



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 10:48 am BST

PDB ID : 4V5A
Title : Structure of the Ribosome Recycling Factor bound to the Thermus thermophilus 70S ribosome with mRNA, ASL-Phe and tRNA-fMet
Authors : Weixlbaumer, A.; Petry, S.; Dunham, C.M.; Selmer, M.; Kelley, A.C.; Ramakrishnan, V.
Deposited on : 2007-06-28
Resolution : 3.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

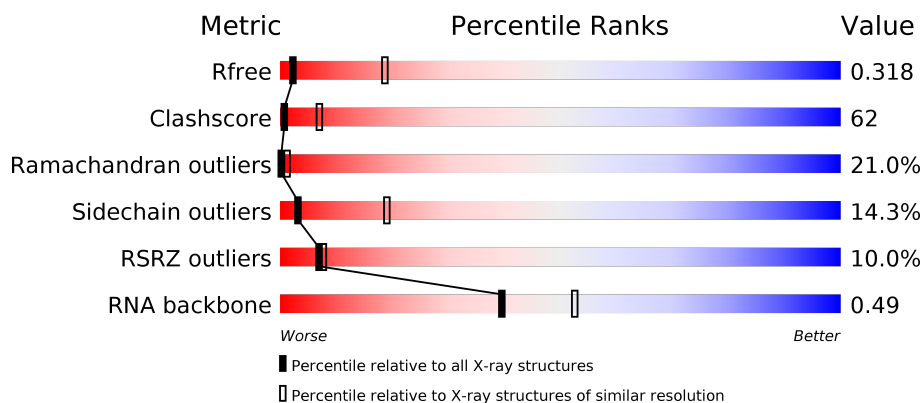
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)
RNA backbone	3102	1002 (4.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AA	1522	<div> <div>11%</div> <div>13% 73% 12% ..</div> </div>
1	CA	1522	<div> <div>13%</div> <div>13% 73% 13% .</div> </div>
2	AB	256	<div> <div>9%</div> <div>16% 50% 25% . 8%</div> </div>
2	CB	256	<div> <div>8%</div> <div>16% 50% 25% . 8%</div> </div>

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Mol	Chain	Length	Quality of chain
3	AC	239	
3	CC	239	
4	AD	209	
4	CD	209	
5	AE	162	
5	CE	162	
6	AF	101	
6	CF	101	
7	AG	156	
7	CG	156	
8	AH	138	
8	CH	138	
9	AI	128	
9	CI	128	
10	AJ	105	
10	CJ	105	
11	AK	129	
11	CK	129	
12	AL	135	
12	CL	135	
13	AM	126	
13	CM	126	
14	AN	61	
14	CN	61	
15	AO	89	

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Mol	Chain	Length	Quality of chain
15	CO	89	
16	AP	88	
16	CP	88	
17	AQ	105	
17	CQ	105	
18	AR	88	
18	CR	88	
19	AS	93	
19	CS	93	
20	AT	106	
20	CT	106	
21	AU	27	
21	CU	27	
22	AV	76	
22	CV	76	
23	AW	77	
23	CW	77	
24	AX	31	
24	CX	31	
25	AY	185	
25	CY	185	
26	B0	85	
26	D0	85	
27	B1	98	
27	D1	98	

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Mol	Chain	Length	Quality of chain
28	B2	72	
28	D2	72	
29	B3	60	
29	D3	60	
30	B4	71	
30	D4	71	
31	B5	60	
31	D5	60	
32	B6	54	
32	D6	54	
33	B7	49	
33	D7	49	
34	B8	65	
34	D8	65	
35	BA	2782	
35	DA	2782	
36	BB	122	
36	DB	122	
37	BC	229	
37	DC	229	
38	BD	276	
38	DD	276	
39	BE	206	
39	DE	206	
40	BF	210	

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Mol	Chain	Length	Quality of chain
40	DF	210	
41	BG	182	
41	DG	182	
42	BH	180	
42	DH	180	
43	BI	148	
43	DI	148	
44	BN	140	
44	DN	140	
45	BO	122	
45	DO	122	
46	BP	150	
46	DP	150	
47	BQ	141	
47	DQ	141	
48	BR	118	
48	DR	118	
49	BS	112	
49	DS	112	
50	BT	146	
50	DT	146	
51	BU	118	
51	DU	118	
52	BV	101	
52	DV	101	

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Mol	Chain	Length	Quality of chain
53	BW	113	
53	DW	113	
54	BX	96	
54	DX	96	
55	BY	110	
55	DY	110	
56	BZ	206	
56	DZ	206	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
23	5MU	CW	55	-	-	-	X
57	MG	AA	1615	-	-	-	X
57	MG	AA	1650	-	-	-	X
57	MG	AA	1653	-	-	-	X
57	MG	AA	1706	-	-	-	X
57	MG	AA	1720	-	-	-	X
57	MG	AA	1733	-	-	-	X
57	MG	AA	1734	-	-	-	X
57	MG	AA	1743	-	-	-	X
57	MG	AA	1749	-	-	-	X
57	MG	AA	1762	-	-	-	X
57	MG	AA	1769	-	-	-	X
57	MG	AA	1780	-	-	-	X
57	MG	AA	1789	-	-	-	X
57	MG	AA	1791	-	-	-	X
57	MG	AA	1798	-	-	-	X
57	MG	AK	201	-	-	-	X
57	MG	AV	102	-	-	-	X
57	MG	AW	108	-	-	-	X
57	MG	AW	123	-	-	-	X
57	MG	BA	3006	-	-	-	X
57	MG	BA	3014	-	-	-	X
57	MG	BA	3035	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
57	MG	BA	3101	-	-	-	X
57	MG	BA	3189	-	-	-	X
57	MG	BA	3207	-	-	-	X
57	MG	BA	3227	-	-	-	X
57	MG	BA	3231	-	-	-	X
57	MG	BA	3251	-	-	-	X
57	MG	BA	3281	-	-	-	X
57	MG	BA	3286	-	-	-	X
57	MG	BA	3287	-	-	-	X
57	MG	BA	3292	-	-	-	X
57	MG	BA	3299	-	-	-	X
57	MG	BA	3309	-	-	-	X
57	MG	BA	3316	-	-	-	X
57	MG	BA	3322	-	-	-	X
57	MG	BA	3334	-	-	-	X
57	MG	BA	3379	-	-	-	X
57	MG	BA	3397	-	-	-	X
57	MG	BA	3412	-	-	-	X
57	MG	BA	3423	-	-	-	X
57	MG	BB	204	-	-	-	X
57	MG	BB	209	-	-	-	X
57	MG	BB	210	-	-	-	X
57	MG	BB	213	-	-	-	X
57	MG	BB	220	-	-	-	X
57	MG	CA	1602	-	-	-	X
57	MG	CA	1643	-	-	-	X
57	MG	CA	1646	-	-	-	X
57	MG	CA	1658	-	-	-	X
57	MG	CA	1681	-	-	-	X
57	MG	CA	1686	-	-	-	X
57	MG	CA	1718	-	-	-	X
57	MG	CA	1725	-	-	-	X
57	MG	CA	1731	-	-	-	X
57	MG	CA	1736	-	-	-	X
57	MG	CA	1737	-	-	-	X
57	MG	CA	1738	-	-	-	X
57	MG	CA	1746	-	-	-	X
57	MG	CA	1747	-	-	-	X
57	MG	CA	1759	-	-	-	X
57	MG	CA	1760	-	-	-	X
57	MG	CA	1765	-	-	-	X
57	MG	CA	1767	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
57	MG	CA	1770	-	-	-	X
57	MG	CA	1785	-	-	-	X
57	MG	CA	1786	-	-	-	X
57	MG	CA	1797	-	-	-	X
57	MG	CE	202	-	-	-	X
57	MG	CN	101	-	-	-	X
57	MG	CV	102	-	-	-	X
57	MG	CW	106	-	-	-	X
57	MG	CW	110	-	-	-	X
57	MG	CW	111	-	-	-	X
57	MG	CW	112	-	-	-	X
57	MG	CW	113	-	-	-	X
57	MG	CW	115	-	-	-	X
57	MG	CW	116	-	-	-	X
57	MG	CW	118	-	-	-	X
57	MG	D1	104	-	-	-	X
57	MG	D2	2601	-	-	-	X
57	MG	DA	3004	-	-	-	X
57	MG	DA	3006	-	-	-	X
57	MG	DA	3049	-	-	-	X
57	MG	DA	3090	-	-	-	X
57	MG	DA	3101	-	-	-	X
57	MG	DA	3127	-	-	-	X
57	MG	DA	3135	-	-	-	X
57	MG	DA	3139	-	-	-	X
57	MG	DA	3140	-	-	-	X
57	MG	DA	3152	-	-	-	X
57	MG	DA	3174	-	-	-	X
57	MG	DA	3176	-	-	-	X
57	MG	DA	3212	-	-	-	X
57	MG	DA	3215	-	-	-	X
57	MG	DA	3220	-	-	-	X
57	MG	DA	3221	-	-	-	X
57	MG	DA	3241	-	-	-	X
57	MG	DA	3258	-	-	-	X
57	MG	DA	3262	-	-	-	X
57	MG	DA	3270	-	-	-	X
57	MG	DA	3271	-	-	-	X
57	MG	DA	3273	-	-	-	X
57	MG	DA	3281	-	-	-	X
57	MG	DA	3282	-	-	-	X
57	MG	DA	3296	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
57	MG	DA	3299	-	-	-	X
57	MG	DA	3301	-	-	-	X
57	MG	DA	3303	-	-	-	X
57	MG	DA	3307	-	-	-	X
57	MG	DA	3326	-	-	-	X
57	MG	DA	3329	-	-	-	X
57	MG	DA	3346	-	-	-	X
57	MG	DA	3358	-	-	-	X
57	MG	DA	3361	-	-	-	X
57	MG	DA	3363	-	-	-	X
57	MG	DA	3367	-	-	-	X
57	MG	DA	3370	-	-	-	X
57	MG	DA	3400	-	-	-	X
57	MG	DA	3409	-	-	-	X
57	MG	DA	3436	-	-	-	X
57	MG	DA	3439	-	-	-	X
57	MG	DA	3441	-	-	-	X
57	MG	DB	202	-	-	-	X
57	MG	DB	219	-	-	-	X
57	MG	DH	201	-	-	-	X

2 Entry composition

There are 58 unique types of molecules in this entry. The entry contains 290487 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	1504	Total	C	N	O	P	0	0	0
			32329	14390	5992	10444	1503			
1	CA	1504	Total	C	N	O	P	0	0	0
			32329	14390	5992	10444	1503			

- Molecule 2 is a protein called 30S RIBOSOMAL PROTEIN S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AB	235	Total	C	N	O	S	0	0	1
			1901	1213	342	341	5			
2	CB	235	Total	C	N	O	S	0	0	1
			1901	1213	342	341	5			

- Molecule 3 is a protein called 30S RIBOSOMAL PROTEIN S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AC	207	Total	C	N	O	S	0	0	1
			1613	1016	315	281	1			
3	CC	207	Total	C	N	O	S	0	0	1
			1613	1016	315	281	1			

- Molecule 4 is a protein called 30S RIBOSOMAL PROTEIN S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AD	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			
4	CD	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

- Molecule 5 is a protein called 30S RIBOSOMAL PROTEIN S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AE	151	Total	C	N	O	S	0	0	1
			1147	724	218	201	4			
5	CE	151	Total	C	N	O	S	0	0	1
			1147	724	218	201	4			

- Molecule 6 is a protein called 30S RIBOSOMAL PROTEIN S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AF	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			
6	CF	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

- Molecule 7 is a protein called 30S RIBOSOMAL PROTEIN S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AG	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			
7	CG	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

- Molecule 8 is a protein called 30S RIBOSOMAL PROTEIN S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AH	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			
8	CH	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

- Molecule 9 is a protein called 30S RIBOSOMAL PROTEIN S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	AI	127	Total	C	N	O	0	0	0
			1011	639	198	174			
9	CI	127	Total	C	N	O	0	0	0
			1011	639	198	174			

- Molecule 10 is a protein called 30S RIBOSOMAL PROTEIN S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AJ	99	Total	C	N	O	S	0	0	1
			795	499	157	138	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	CJ	99	Total	C	N	O	S	0	0	1
			795	499	157	138	1			

- Molecule 11 is a protein called 30S RIBOSOMAL PROTEIN S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AK	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			
11	CK	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			

- Molecule 12 is a protein called 30S RIBOSOMAL PROTEIN S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AL	125	Total	C	N	O	S	0	0	1
			971	611	196	163	1			
12	CL	125	Total	C	N	O	S	0	0	1
			971	611	196	163	1			

- Molecule 13 is a protein called 30S RIBOSOMAL PROTEIN S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AM	125	Total	C	N	O	S	0	0	1
			988	611	206	169	2			
13	CM	125	Total	C	N	O	S	0	0	1
			988	611	206	169	2			

- Molecule 14 is a protein called 30S RIBOSOMAL PROTEIN S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			
14	CN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

- Molecule 15 is a protein called 30S RIBOSOMAL PROTEIN S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AO	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			
15	CO	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

- Molecule 16 is a protein called 30S RIBOSOMAL PROTEIN S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AP	84	Total	C	N	O	S	0	0	1
			701	443	140	117	1			
16	CP	84	Total	C	N	O	S	0	0	1
			701	443	140	117	1			

- Molecule 17 is a protein called 30S RIBOSOMAL PROTEIN S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AQ	100	Total	C	N	O	S	0	0	1
			824	528	152	142	2			
17	CQ	100	Total	C	N	O	S	0	0	1
			824	528	152	142	2			

- Molecule 18 is a protein called 30S RIBOSOMAL PROTEIN S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	AR	70	Total	C	N	O	0	0	0
			574	367	112	95			
18	CR	70	Total	C	N	O	0	0	0
			574	367	112	95			

- Molecule 19 is a protein called 30S RIBOSOMAL PROTEIN S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AS	79	Total	C	N	O	S	0	0	1
			630	403	115	110	2			
19	CS	79	Total	C	N	O	S	0	0	1
			630	403	115	110	2			

- Molecule 20 is a protein called 30S RIBOSOMAL PROTEIN S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AT	99	Total	C	N	O	S	0	0	0
			763	470	162	129	2			
20	CT	99	Total	C	N	O	S	0	0	0
			763	470	162	129	2			

- Molecule 21 is a protein called 30S RIBOSOMAL PROTEIN THX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	AU	25	Total	C	N	O	0	0	1
			209	128	51	30			
21	CU	25	Total	C	N	O	0	0	1
			209	128	51	30			

- Molecule 22 is a RNA chain called P-SITE RNA ASL-PHE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	AV	17	Total	C	N	O	P	0	0	0
			362	163	68	115	16			
22	CV	17	Total	C	N	O	P	0	0	0
			362	163	68	115	16			

- Molecule 23 is a RNA chain called TRNA-FMET.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	AW	77	Total	C	N	O	P	0	0	0
			1641	733	297	535	76			
23	CW	77	Total	C	N	O	P	0	0	0
			1641	733	297	535	76			

- Molecule 24 is a RNA chain called MRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	AX	11	Total	C	N	O	P	0	0	0
			230	104	38	78	10			
24	CX	11	Total	C	N	O	P	0	0	0
			230	104	38	78	10			

- Molecule 25 is a protein called RIBOSOME RECYCLING FACTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	AY	185	Total	C	N	O	S	0	0	0
			1478	924	270	282	2			
25	CY	185	Total	C	N	O	S	0	0	0
			1478	924	270	282	2			

- Molecule 26 is a protein called 50S RIBOSOMAL PROTEIN L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	B0	85	Total	C	N	O	S	0	0	0
			652	403	137	111	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	D0	85	Total	C	N	O	S	0	0	0
			652	403	137	111	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B0	6	ALA	GLY	conflict	UNP P60493
B0	8	ALA	GLY	conflict	UNP P60493
D0	6	ALA	GLY	conflict	UNP P60493
D0	8	ALA	GLY	conflict	UNP P60493

- Molecule 27 is a protein called 50S RIBOSOMAL PROTEIN L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	B1	89	Total	C	N	O		0	0	1
			693	435	140	118				
27	D1	89	Total	C	N	O		0	0	1
			693	435	140	118				

- Molecule 28 is a protein called 50S RIBOSOMAL PROTEIN L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	B2	51	Total	C	N	O	S	0	0	1
			421	263	85	72	1			
28	D2	51	Total	C	N	O	S	0	0	1
			421	263	85	72	1			

- Molecule 29 is a protein called 50S RIBOSOMAL PROTEIN L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	B3	60	Total	C	N	O	S	0	0	1
			468	298	91	78	1			
29	D3	60	Total	C	N	O	S	0	0	1
			468	298	91	78	1			

- Molecule 30 is a protein called 50S RIBOSOMAL PROTEIN L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	B4	50	Total	C	N	O		0	0	1
			242	143	50	49				
30	D4	50	Total	C	N	O		0	0	1
			242	143	50	49				

- Molecule 31 is a protein called 50S RIBOSOMAL PROTEIN L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	B5	59	Total	C	N	O	S	0	0	0
			459	288	90	76	5			
31	D5	59	Total	C	N	O	S	0	0	0
			459	288	90	76	5			

- Molecule 32 is a protein called 50S RIBOSOMAL PROTEIN L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	B6	45	Total	C	N	O	S	0	0	1
			381	235	78	64	4			
32	D6	45	Total	C	N	O	S	0	0	1
			381	235	78	64	4			

- Molecule 33 is a protein called 50S RIBOSOMAL PROTEIN L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	B7	49	Total	C	N	O	S	0	0	1
			419	257	105	55	2			
33	D7	49	Total	C	N	O	S	0	0	1
			419	257	105	55	2			

- Molecule 34 is a protein called 50S RIBOSOMAL PROTEIN L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	B8	64	Total	C	N	O	S	0	0	1
			508	326	102	78	2			
34	D8	64	Total	C	N	O	S	0	0	1
			508	326	102	78	2			

- Molecule 35 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BA	2767	Total	C	N	O	P	0	0	0
			59601	26526	11153	19156	2766			
35	DA	2767	Total	C	N	O	P	0	0	0
			59601	26526	11153	19156	2766			

- Molecule 36 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	BB	119	Total	C	N	O	P	0	0	0
			2551	1136	471	826	118			
36	DB	119	Total	C	N	O	P	0	0	0
			2551	1136	471	826	118			

- Molecule 37 is a protein called 50S RIBOSOMAL PROTEIN L1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
37	BC	191	Total	C	N	O	0	0	1
			1142	691	221	230			
37	DC	191	Total	C	N	O	0	0	1
			1143	692	221	230			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
DC	106	ALA	GLY	conflict	UNP Q5SLP7
DC	132	ALA	GLY	conflict	UNP Q5SLP7

- Molecule 38 is a protein called 50S RIBOSOMAL PROTEIN L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	BD	272	Total	C	N	O	S	0	0	1
			2105	1329	417	356	3			
38	DD	272	Total	C	N	O	S	0	0	1
			2105	1329	417	356	3			

- Molecule 39 is a protein called 50S RIBOSOMAL PROTEIN L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	BE	205	Total	C	N	O	S	0	0	1
			1564	988	300	270	6			
39	DE	205	Total	C	N	O	S	0	0	1
			1564	988	300	270	6			

- Molecule 40 is a protein called 50S RIBOSOMAL PROTEIN L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	BF	208	Total	C	N	O	S	0	0	1
			1624	1035	304	282	3			
40	DF	208	Total	C	N	O	S	0	0	1
			1624	1035	304	282	3			

- Molecule 41 is a protein called 50S RIBOSOMAL PROTEIN L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BG	181	Total	C	N	O	S	0	0	0
			1474	942	268	260	4			
41	DG	181	Total	C	N	O	S	0	0	0
			1474	942	268	260	4			

- Molecule 42 is a protein called 50S RIBOSOMAL PROTEIN L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BH	160	Total	C	N	O	S	0	0	1
			1223	773	229	220	1			
42	DH	160	Total	C	N	O	S	0	0	1
			1223	773	229	220	1			

- Molecule 43 is a protein called 50S RIBOSOMAL PROTEIN L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	BI	146	Total	C	N	O	S	0	0	1
			1132	723	201	207	1			
43	DI	146	Total	C	N	O	S	0	0	1
			1132	723	201	207	1			

- Molecule 44 is a protein called 50S RIBOSOMAL PROTEIN L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	BN	139	Total	C	N	O	S	0	0	1
			1105	712	207	182	4			
44	DN	139	Total	C	N	O	S	0	0	1
			1105	712	207	182	4			

- Molecule 45 is a protein called 50S RIBOSOMAL PROTEIN L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	BO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			
45	DO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			

- Molecule 46 is a protein called 50S RIBOSOMAL PROTEIN L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	BP	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			
46	DP	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			

- Molecule 47 is a protein called 50S RIBOSOMAL PROTEIN L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	BQ	136	Total	C	N	O	S	0	0	0
			1080	688	204	183	5			
47	DQ	136	Total	C	N	O	S	0	0	0
			1080	688	204	183	5			

- Molecule 48 is a protein called 50S RIBOSOMAL PROTEIN L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	BR	117	Total	C	N	O		0	0	0
			960	599	202	159				
48	DR	117	Total	C	N	O		0	0	0
			960	599	202	159				

- Molecule 49 is a protein called 50S RIBOSOMAL PROTEIN L18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	BS	99	Total	C	N	O		0	0	1
			771	486	155	130				
49	DS	99	Total	C	N	O		0	0	1
			771	486	155	130				

- Molecule 50 is a protein called 50S RIBOSOMAL PROTEIN L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	BT	138	Total	C	N	O	S	0	0	1
			1142	710	235	196	1			
50	DT	138	Total	C	N	O	S	0	0	1
			1142	710	235	196	1			

- Molecule 51 is a protein called 50S RIBOSOMAL PROTEIN L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	BU	117	Total	C	N	O	S	0	0	0
			958	604	202	151	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	DU	117	Total	C	N	O	S	0	0	0
			958	604	202	151	1			

- Molecule 52 is a protein called 50S RIBOSOMAL PROTEIN L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	BV	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			
52	DV	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			

- Molecule 53 is a protein called 50S RIBOSOMAL PROTEIN L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	BW	113	Total	C	N	O	S	0	0	0
			896	563	176	155	2			
53	DW	113	Total	C	N	O	S	0	0	0
			896	563	176	155	2			

- Molecule 54 is a protein called 50S RIBOSOMAL PROTEIN L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	BX	93	Total	C	N	O		0	0	1
			726	471	132	123				
54	DX	93	Total	C	N	O		0	0	1
			726	471	132	123				

- Molecule 55 is a protein called 50S RIBOSOMAL PROTEIN L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	BY	101	Total	C	N	O	S	0	0	1
			776	500	149	123	4			
55	DY	101	Total	C	N	O	S	0	0	1
			776	500	149	123	4			

- Molecule 56 is a protein called 50S RIBOSOMAL PROTEIN L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
56	BZ	177	Total	C	N	O	S	0	0	1
			1404	897	253	252	2			
56	DZ	177	Total	C	N	O	S	0	0	1
			1404	897	253	252	2			

- Molecule 57 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	BA	445	Total 445	Mg 445	0	0
57	AK	1	Total 1	Mg 1	0	0
57	DF	5	Total 5	Mg 5	0	0
57	CV	4	Total 4	Mg 4	0	0
57	D2	2	Total 2	Mg 2	0	0
57	BE	1	Total 1	Mg 1	0	0
57	AW	23	Total 23	Mg 23	0	0
57	DU	3	Total 3	Mg 3	0	0
57	B1	4	Total 4	Mg 4	0	0
57	BP	3	Total 3	Mg 3	0	0
57	AX	5	Total 5	Mg 5	0	0
57	CN	1	Total 1	Mg 1	0	0
57	DN	2	Total 2	Mg 2	0	0
57	CA	208	Total 208	Mg 208	0	0
57	B5	2	Total 2	Mg 2	0	0
57	BB	20	Total 20	Mg 20	0	0
57	AE	2	Total 2	Mg 2	0	0
57	DG	1	Total 1	Mg 1	0	0
57	D3	1	Total 1	Mg 1	0	0
57	BF	4	Total 4	Mg 4	0	0
57	AV	4	Total 4	Mg 4	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	BX	2	Total 2	Mg 2	0	0
57	B2	2	Total 2	Mg 2	0	0
57	AA	204	Total 204	Mg 204	0	0
57	D7	1	Total 1	Mg 1	0	0
57	CX	4	Total 4	Mg 4	0	0
57	DV	1	Total 1	Mg 1	0	0
57	AM	1	Total 1	Mg 1	0	0
57	BU	2	Total 2	Mg 2	0	0
57	AD	2	Total 2	Mg 2	0	0
57	BN	4	Total 4	Mg 4	0	0
57	DH	1	Total 1	Mg 1	0	0
57	BG	1	Total 1	Mg 1	0	0
57	DS	1	Total 1	Mg 1	0	0
57	DE	2	Total 2	Mg 2	0	0
57	B3	1	Total 1	Mg 1	0	0
57	DX	3	Total 3	Mg 3	0	0
57	DA	441	Total 441	Mg 441	0	0
57	B7	1	Total 1	Mg 1	0	0
57	AL	2	Total 2	Mg 2	0	0
57	AG	1	Total 1	Mg 1	0	0
57	BO	1	Total 1	Mg 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	D1	4	Total 4	Mg 4	0	0
57	DP	4	Total 4	Mg 4	0	0
57	CW	23	Total 23	Mg 23	0	0
57	D5	1	Total 1	Mg 1	0	0
57	BD	3	Total 3	Mg 3	0	0
57	CE	3	Total 3	Mg 3	0	0
57	CG	1	Total 1	Mg 1	0	0
57	DD	3	Total 3	Mg 3	0	0
57	CL	2	Total 2	Mg 2	0	0
57	DB	19	Total 19	Mg 19	0	0

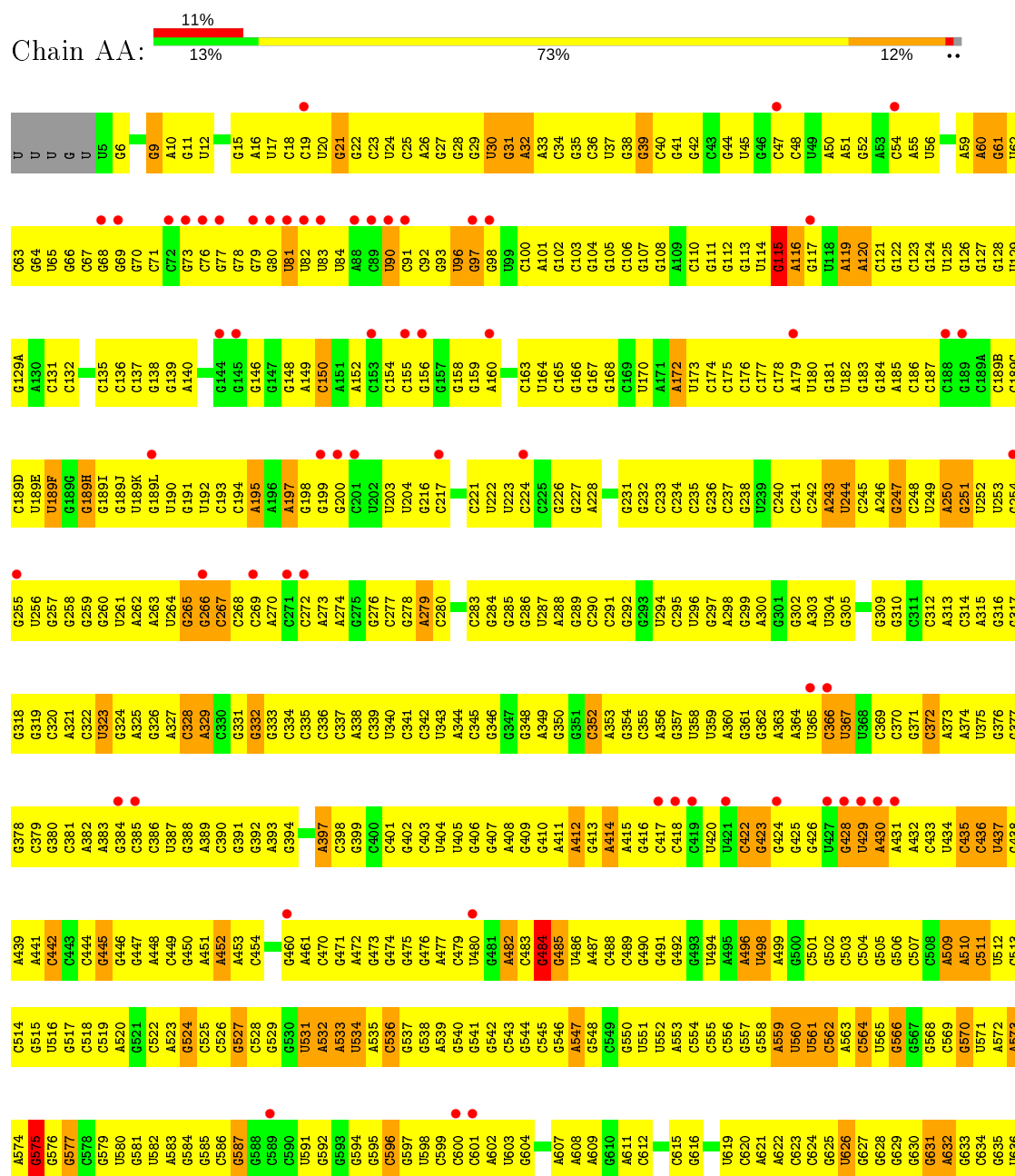
- Molecule 58 is ZINC ION (three-letter code: ZN) (formula: Zn).

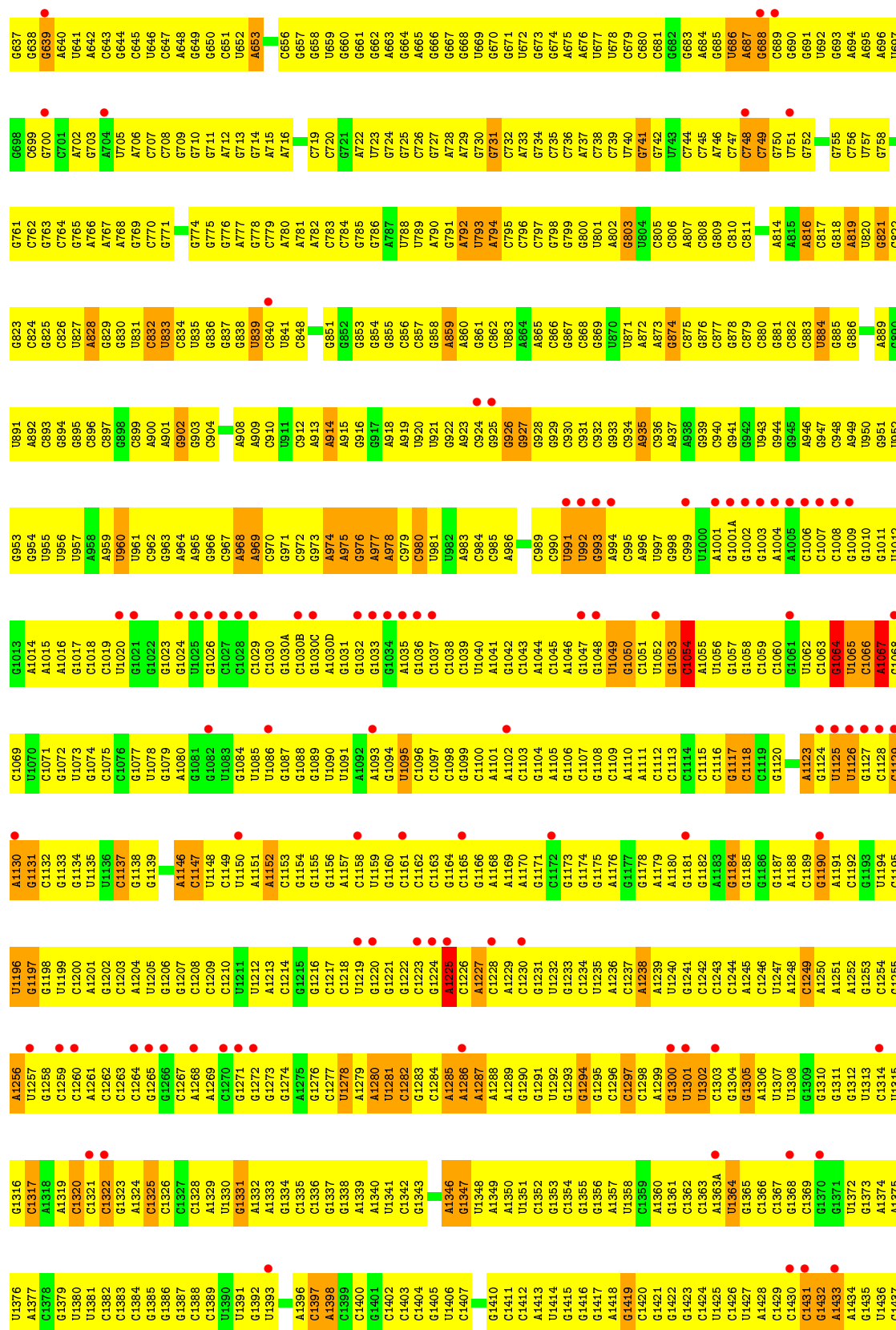
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	CN	1	Total 1	Zn 1	0	0
58	AD	1	Total 1	Zn 1	0	0
58	CD	1	Total 1	Zn 1	0	0
58	AN	1	Total 1	Zn 1	0	0

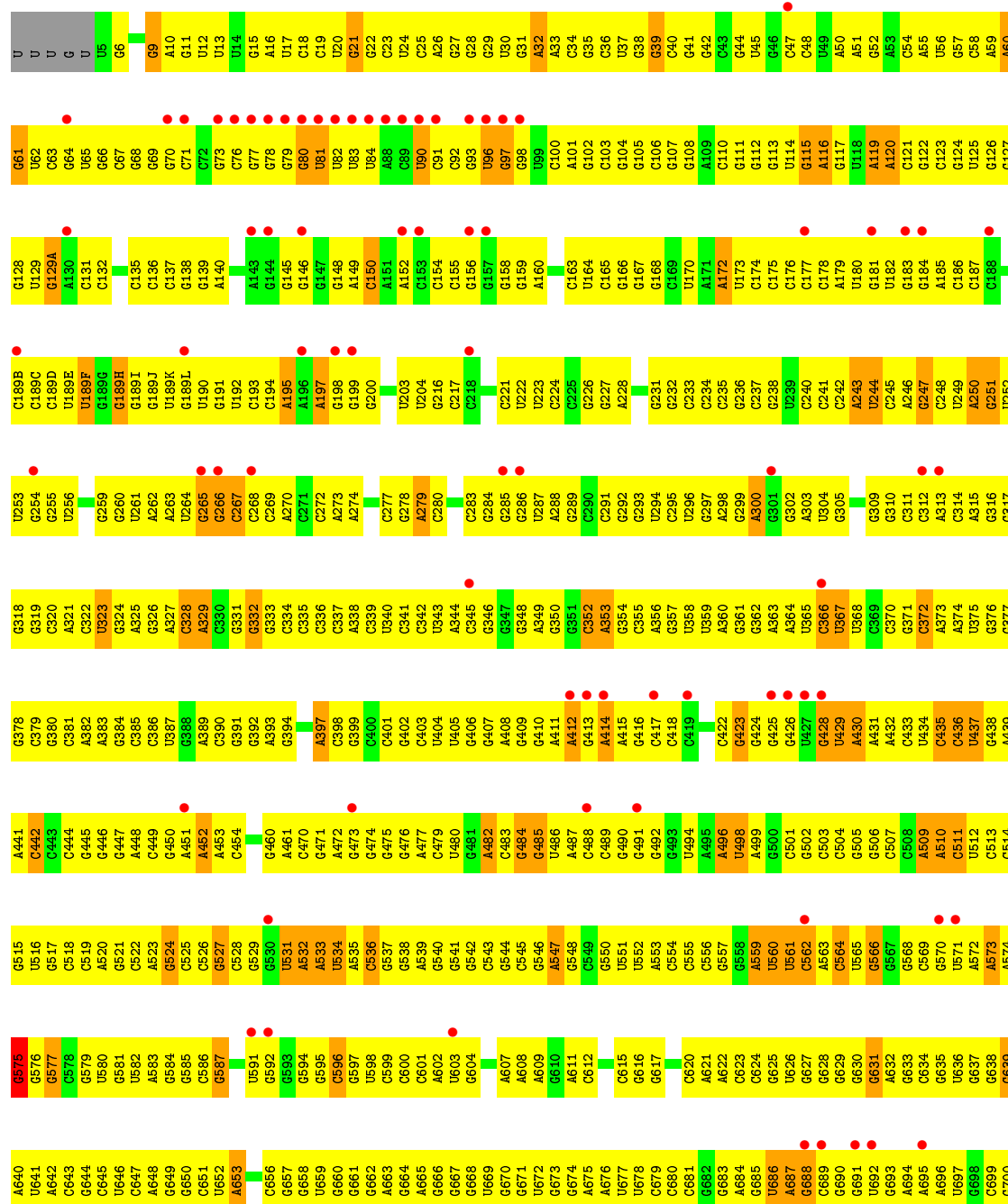
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

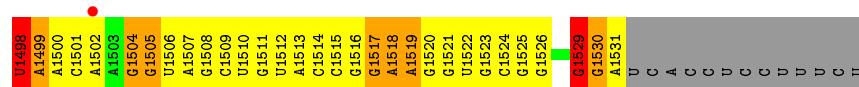
- Molecule 1: 16S ribosomal RNA



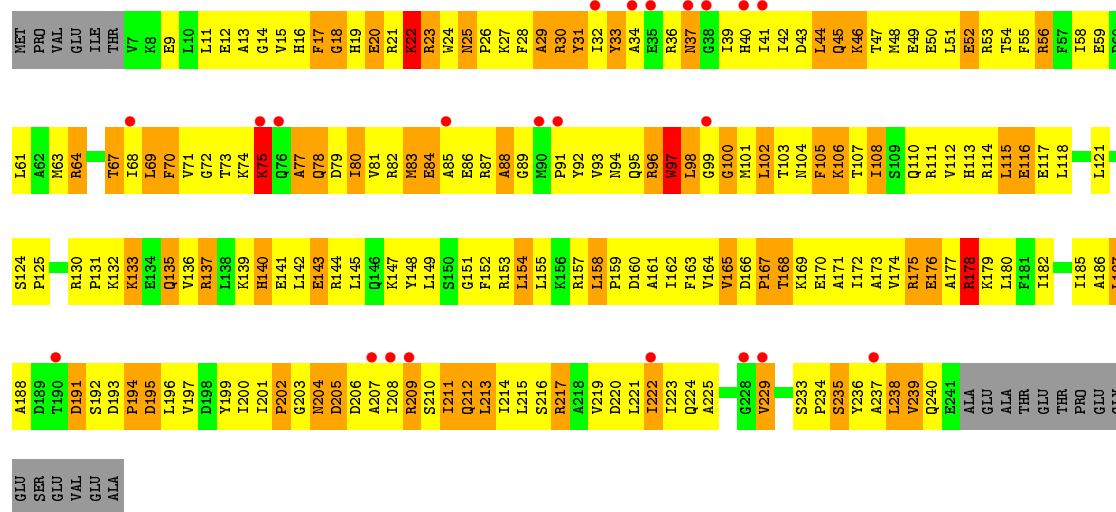
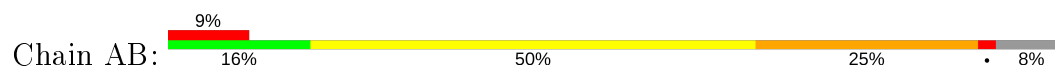




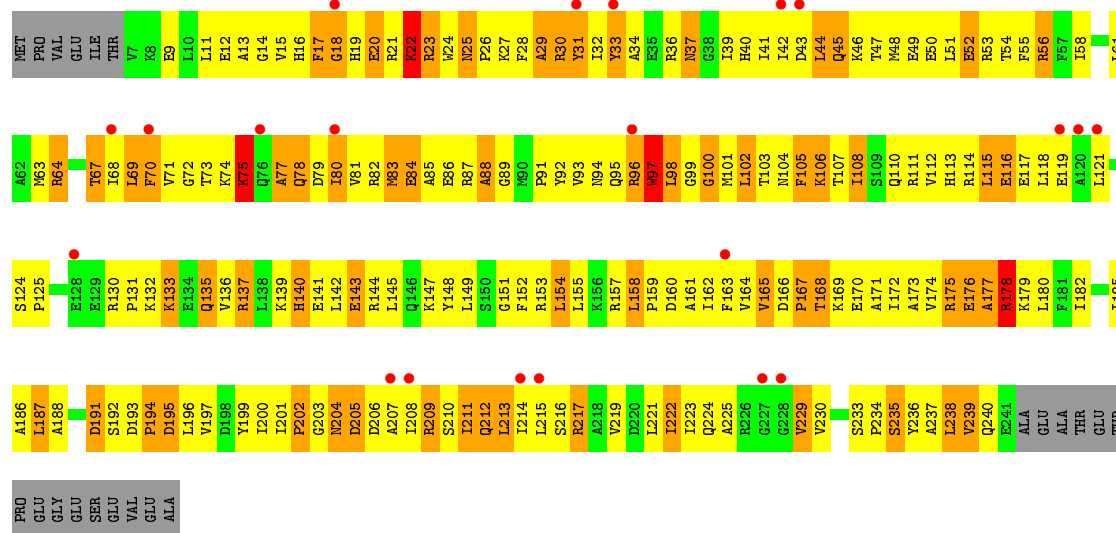
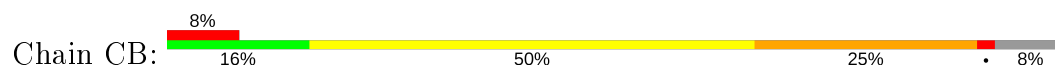
A1433	G1373	U1313	G1253	G1193	C1128	G1068	U1012	U952	A892	G825	G763	C701
A1434	A1374	C1314	C1254	U1194	C1129	C1069	G1013	G953	C893	C826	C764	A702
G1435	A1375	U1315	G1255	U1195	A1130	U1070	A1014	G954	G894	C825	G765	G703
U1436	U1376	G1316	A1256	U1196	G1131	U1071	A1015	U955	G895	A828	A766	A704
G1437	A1377	G1317	G1257	G1197	C1132	U1072	A1016	U956	C896	G829	A767	U705
G1438	G1378	A1318	G1258	U1198	G1133	U1073	G1017	U957	C897	G830	A768	A706
C1439	G1379	A1319	C1259	U1199	G1134	G1074	C1018	A858	G898	U831	C707	C707
G1440	C1380	C1320	G1260	U1200	G1135	G1075	C1019	U959	C899	U832	C708	C708
G1441	A1381	A1321	A1261	A1201	U1136	C1076	U1020	U960	A900	U833	G709	G709
G1442	C1382	C1322	C1262	C1202	C1137	G1077	G1021	U961	A901	C834	G774	G711
G1442A	C1383	C1323	C1263	C1203	G1138	U1078	G1022	G962	G902	U835	G775	G711
A1442B	C1384	A1324	G1264	A1204	G1139	G1079	G1023	G963	G903	G836	A712	A712
G1443	G1385	C1325	G1265	U1205	C1140	A1080	G1024	A964	C904	G837	G776	G713
G1444	G1386	C1326	G1266	G1206	C1141	G1081	U1025	A865	U905	G838	A777	G714
C1445	C1387	G1327	C1267	G1207	A1146	U1082	G1026	G966	G906	U839	A715	A715
U1446	C1388	C1328	A1268	C1208	C1147	U1083	G1027	C967	A907	C840	C779	A716
A1447	C1389	A1329	C1269	C1209	U1084	U1084	C1028	A968	A908	U841	A780	A780
G1456	U1390	U1330	G1270	C1210	U1148	U1085	C1029	A969	A909	C948	A781	C719
G1391	U1391	G1331	G1271	U1211	C1149	U1086	C1030	C970	C910	C849	A782	C720
G1392	U1392	A1332	G1272	U1212	U1150	G1087	G1030A	G971	U911	U850	C783	G721
G1393	U1393	A1333	G1273	A1213	A1151	G1088	G1030B	C972	C912	C851	C784	A722
A1394	A1394	G1334	G1274	C1214	A1152	G1089	G1030C	G973	A913	G852	G785	U723
A1395	A1395	G1335	A1275	G1215	C1153	U1090	A1030D	A974	A914	G853	G786	A723
A1396	A1396	G1336	G1276	G1216	G1154	U1091	G1031	A975	A915	G854	A787	G724
C1397	C1397	G1338	G1277	C1217	G1155	A1092	G1032	G976	G916	G855	G788	G725
A1398	A1398	A1339	U1278	C1218	G1156	A1093	G1033	A977	G917	C856	C726	C726
C1399	C1399	A1340	A1279	U1219	A1157	U1094	U1034	A978	A918	C857	G727	G727
C1400	C1400	U1341	A1280	G1220	C1158	U1095	A1035	C979	A919	C858	A792	A728
G1401	G1401	C1342	U1281	G1221	U1159	C1096	G1036	A859	U920	U793	U793	A729
G1402	C1402	G1343	C1282	G1222	G1160	C1097	C1037	U981	U921	U891	A794	G730
C1403	C1403	C1344	G1283	C1223	C1161	U1098	C1038	U982	G922	G861	C795	G731
C1404	C1404	U1345	C1284	G1224	C1162	G1099	C1039	A883	A823	C862	C732	G732
G1405	G1405	A1346	A1285	A1225	C1163	C1100	U1040	C984	C924	U863	A733	A733
U1471	U1406	G1347	A1286	C1226	G1164	A1101	A1041	C985	G925	C797	G734	G734
U1472	A1407	U1348	A1287	A1227	C1165	A1102	G1042	G986	G926	G798	G735	G735
A1408	A1408	A1349	A1288	C1228	G1166	C1097	G1043	G987	G927	C866	C736	C736
A1409	A1409	A1350	A1289	A1229	A1168	G1104	A1044	C989	G928	G867	A737	A737
G1410	G1410	U1351	G1290	C1230	A1169	A1105	C1045	C990	G929	C868	C738	C738
C1411	C1411	C1352	G1291	G1231	A1170	G1106	A1046	U991	C930	G869	C739	C739
C1412	C1412	G1353	U1292	U1232	C1171	C1107	G1047	U992	C931	U870	U740	U740
C1354	C1354	C1354	G1293	C1233	G1174	G1108	U1048	G993	C932	U871	G741	G741
G1355	G1355	G1355	G1294	C1234	A1175	A1109	U1049	A994	G933	A872	G742	G742
G1356	G1356	C1357	G1295	U1235	A1176	A1110	G1050	C995	C934	A873	U743	U743
A1357	A1357	C1358	C1296	U1236	G1177	C1112	C1051	A996	A835	G874	C744	C744
U1358	U1358	C1359	C1297	C1237	G1178	C1113	U1052	U997	C936	C875	C745	C745
A1418	A1418	C1359	A1298	A1238	G1179	C1114	G1053	G998	A937	C876	A746	A746
C1419	C1419	A1360	A1299	A1239	A1179	C1114	C1054	C999	A938	C877	C747	C747
G1361	G1361	U1361	G1300	U1240	A1180	C1115	U1055	U1000	G939	G878	C748	C748
C1362	C1362	U1301	U1301	G1241	G1181	C1116	U1056	A1001	C940	C879	C749	C749
C1363	C1363	U1302	G1302	C1242	G1182	G1117	G1057	G1001A	G941	C880	U751	U751
A1363A	A1363A	C1303	C1303	C1243	A1183	C1118	G1058	G1002	G942	G881	G752	G752
U1364	U1364	G1304	G1304	C1244	G1184	C1119	G1059	G1003	U943	A816	G755	G755
G1365	G1365	A1245	G1305	A1245	G1185	C1120	C1060	C882	C983	C817	C756	C756
C1366	C1366	A1306	C1306	C1246	G1186	U1121	G1061	G945	G945	U884	U757	U757
G1367	G1367	U1307	U1307	U1247	G1187	U1122	U1062	A946	A946	G885	C758	C758
U1308	U1308	A1248	A1248	C1248	A1188	A1123	C1063	C1007	C948	U820	U757	U757
C1368	C1368	G1309	G1309	C1249	C1189	G1124	G1064	C1008	C948	G886	G758	G758
C1369	C1369	G1310	G1310	A1250	G1190	U1125	U1065	G1009	A949	C821	G761	G761
G1370	G1370	C1311	G1311	A1251	C1191	U1126	U1066	G1010	U950	C822	C762	C762
U1372	U1372	A1252	G1312	A1252	C1192	G1127	A1067	G1011	G951	C824	C762	C762



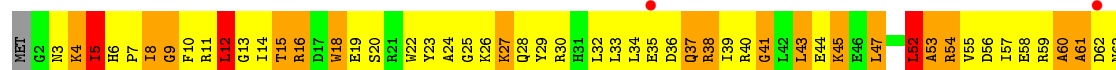
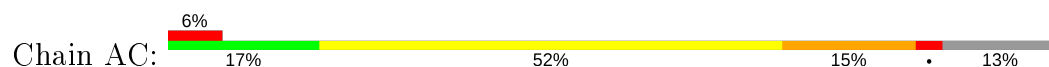
• Molecule 2: 30S RIBOSOMAL PROTEIN S2

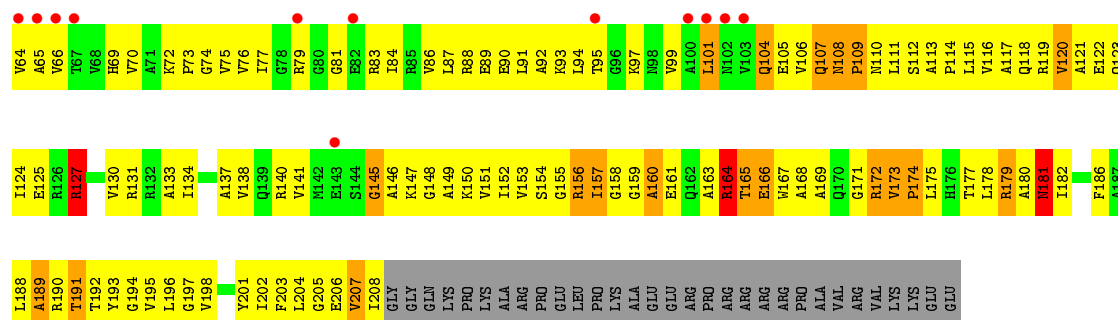


• Molecule 2: 30S RIBOSOMAL PROTEIN S2

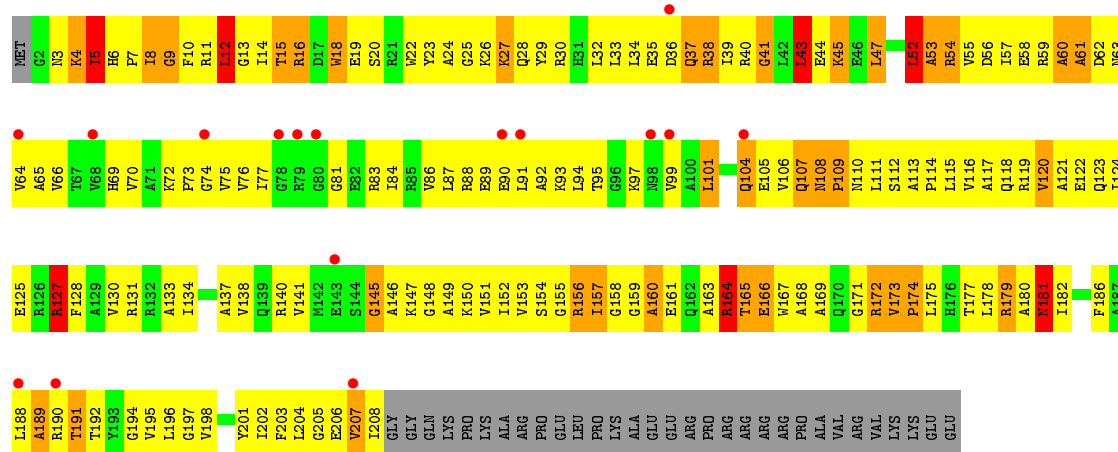
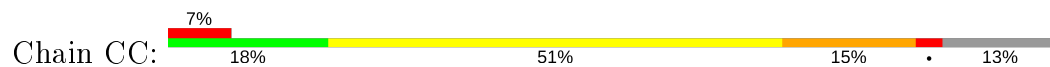


• Molecule 3: 30S RIBOSOMAL PROTEIN S3

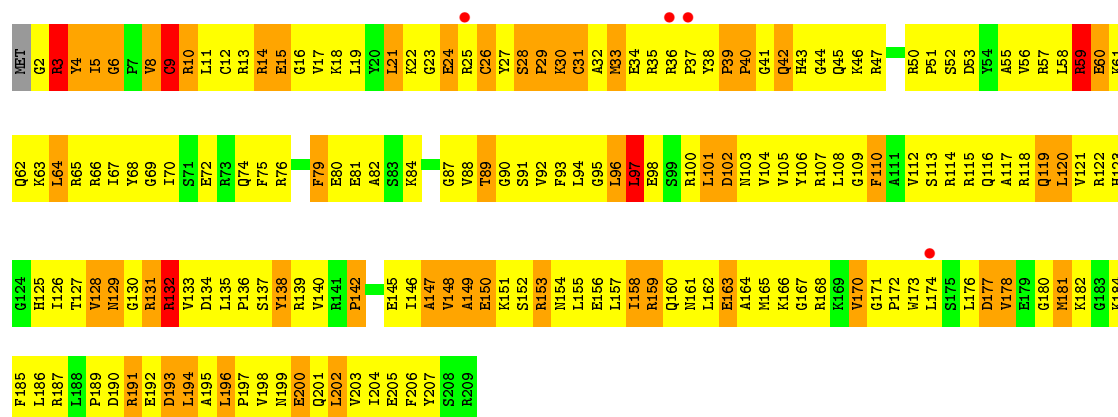
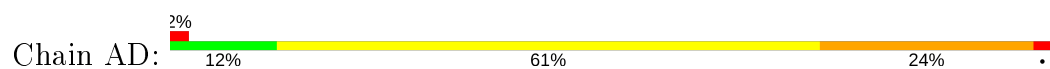




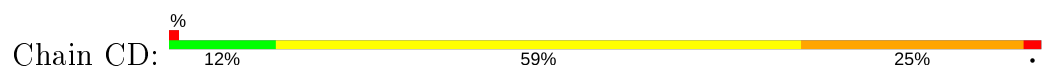
• Molecule 3: 30S RIBOSOMAL PROTEIN S3

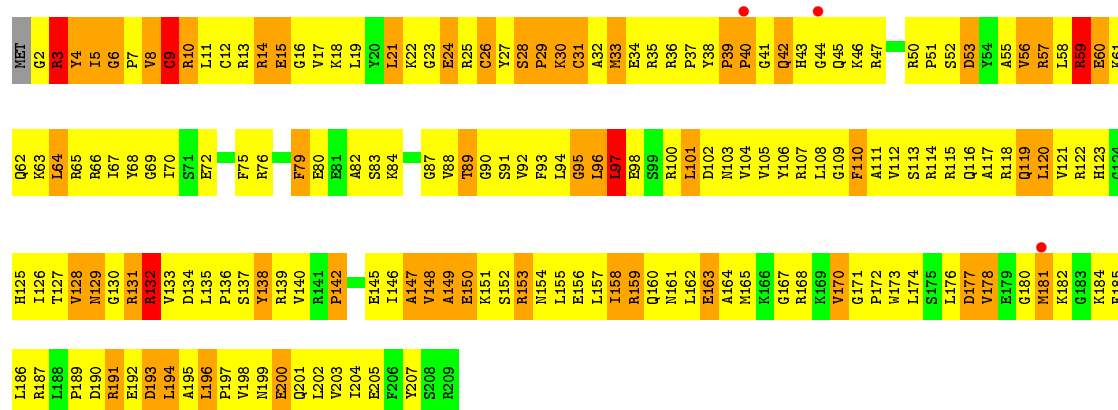


• Molecule 4: 30S RIBOSOMAL PROTEIN S4

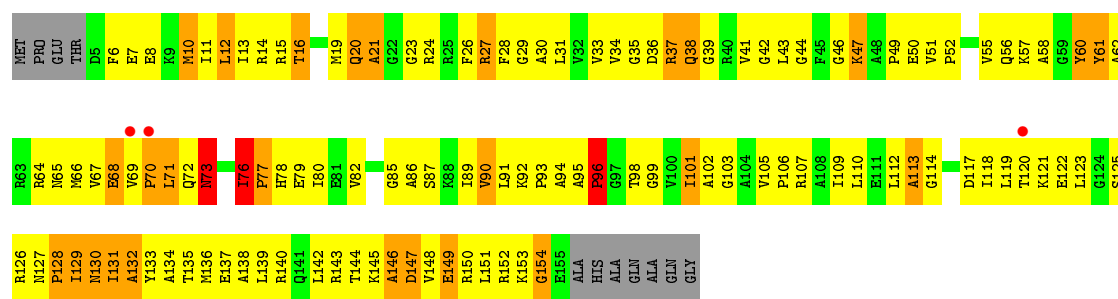


• Molecule 4: 30S RIBOSOMAL PROTEIN S4

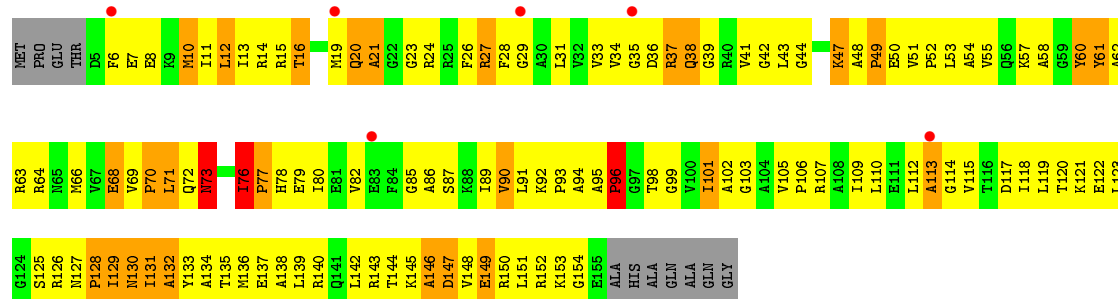




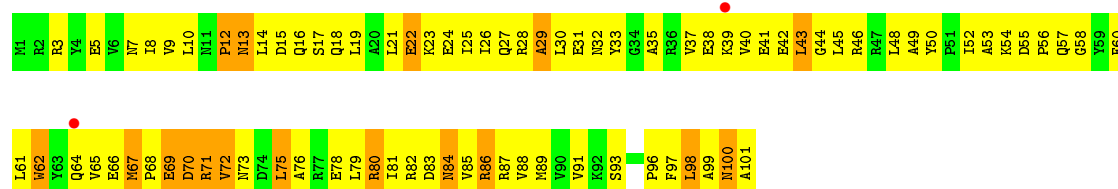
• Molecule 5: 30S RIBOSOMAL PROTEIN S5



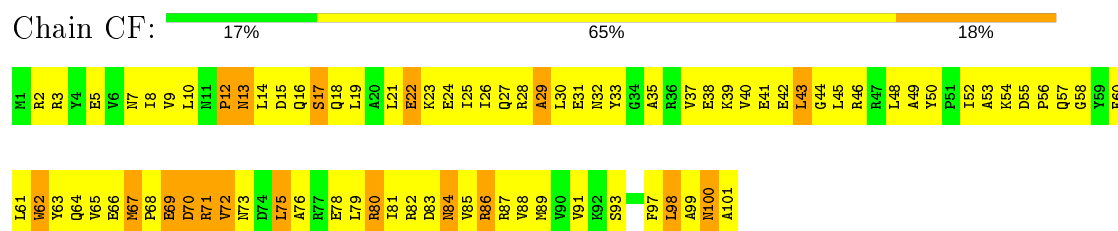
• Molecule 5: 30S RIBOSOMAL PROTEIN S5



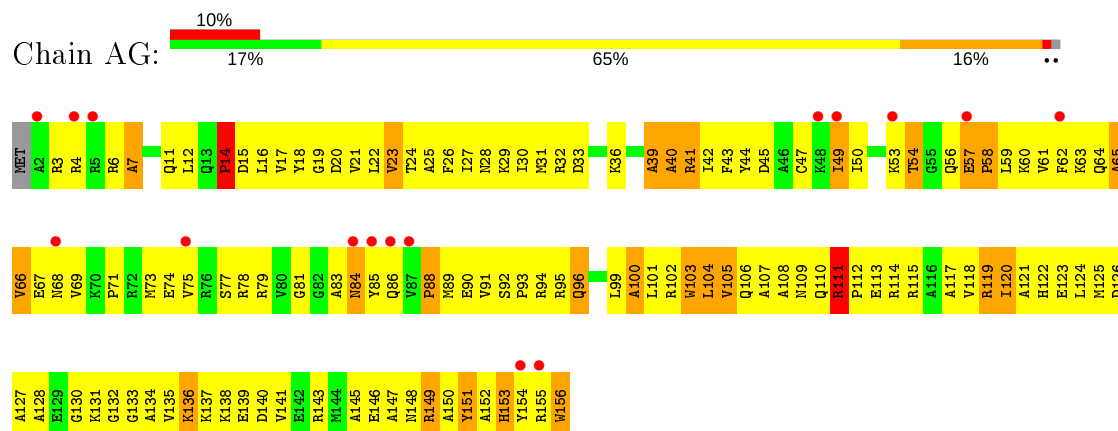
• Molecule 6: 30S RIBOSOMAL PROTEIN S6



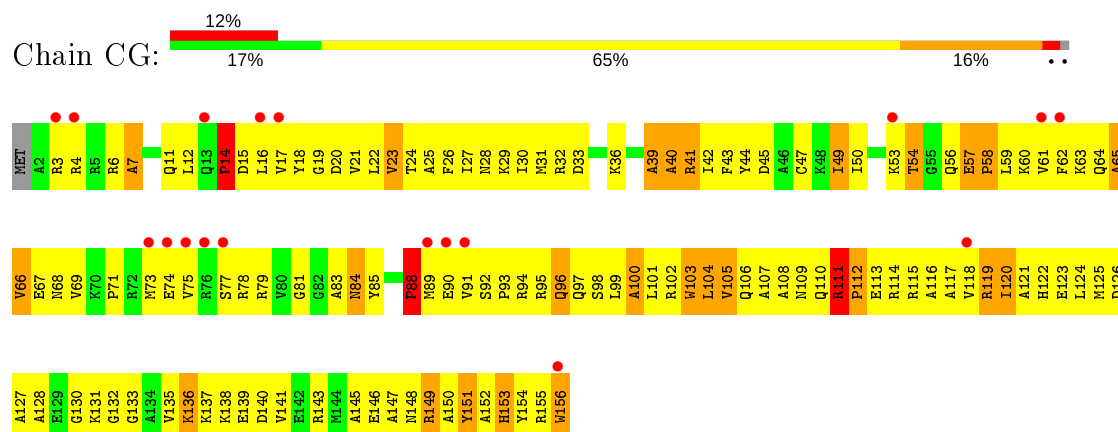
- Molecule 6: 30S RIBOSOMAL PROTEIN S6



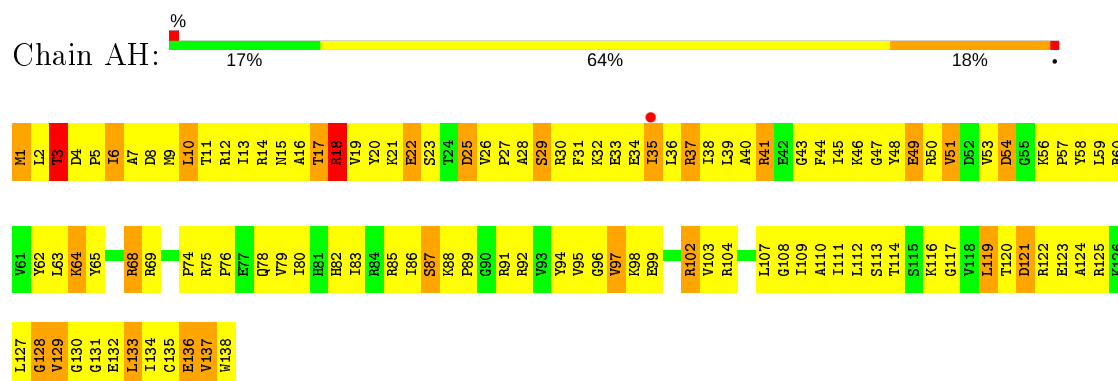
- Molecule 7: 30S RIBOSOMAL PROTEIN S7



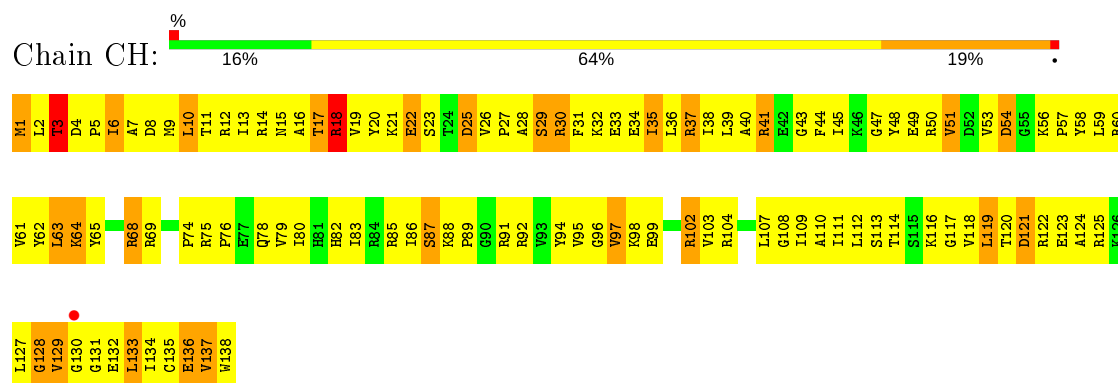
- Molecule 7: 30S RIBOSOMAL PROTEIN S7



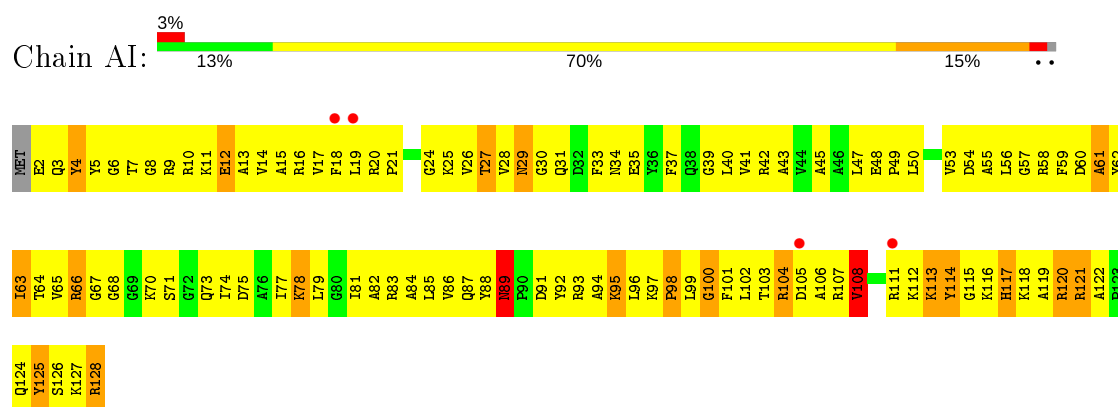
- Molecule 8: 30S RIBOSOMAL PROTEIN S8



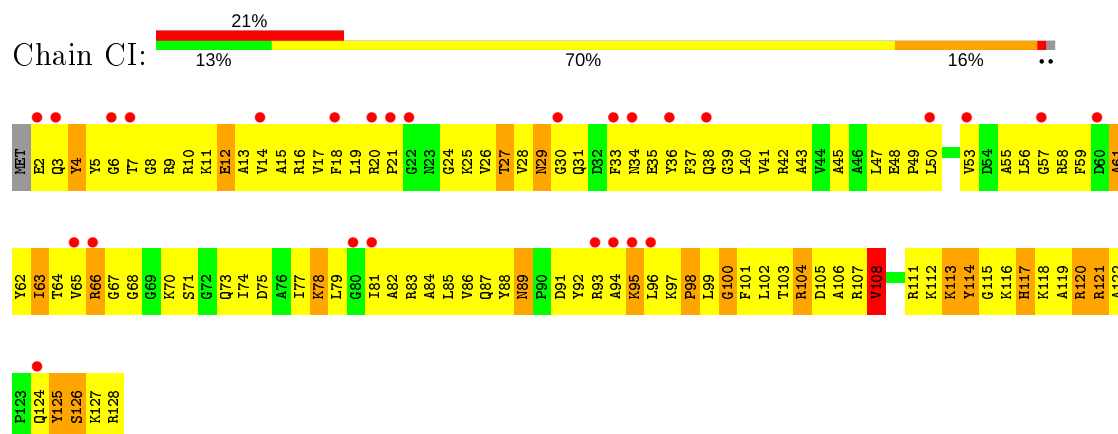
- Molecule 8: 30S RIBOSOMAL PROTEIN S8



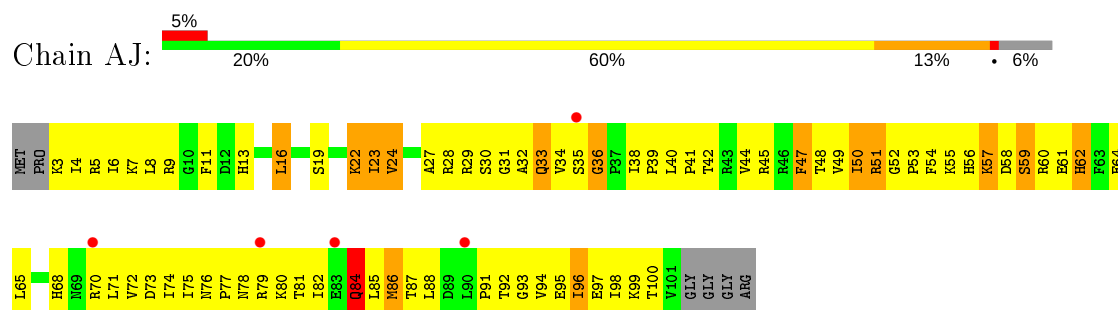
- Molecule 9: 30S RIBOSOMAL PROTEIN S9



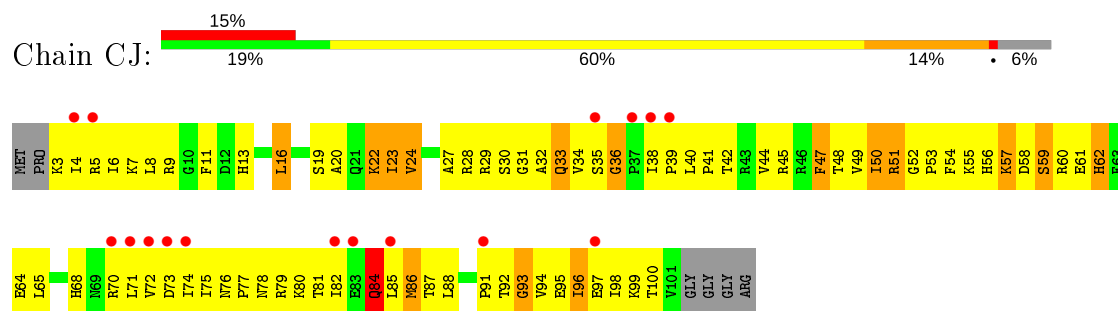
- Molecule 9: 30S RIBOSOMAL PROTEIN S9



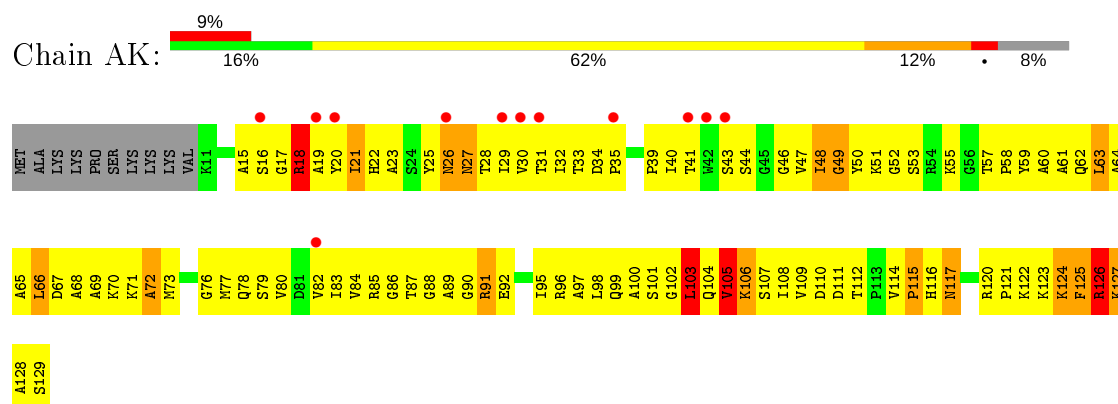
- Molecule 10: 30S RIBOSOMAL PROTEIN S10



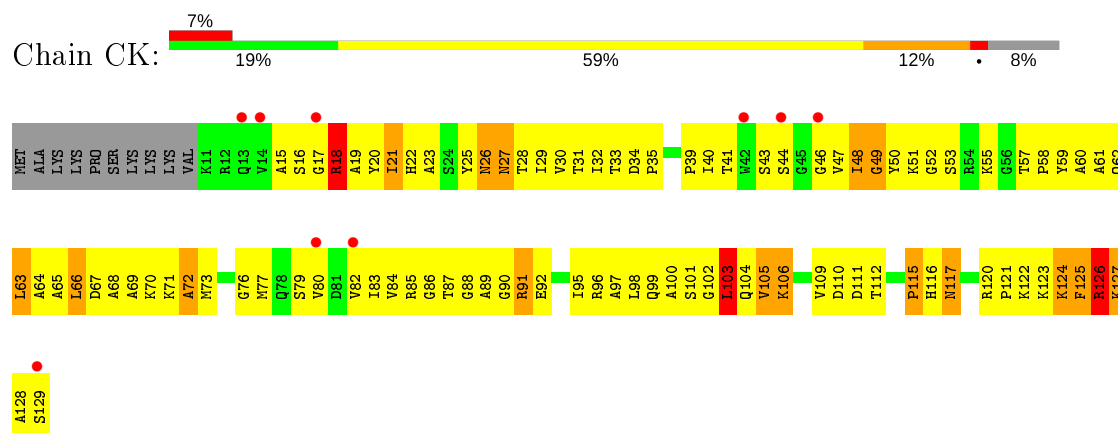
• Molecule 10: 30S RIBOSOMAL PROTEIN S10



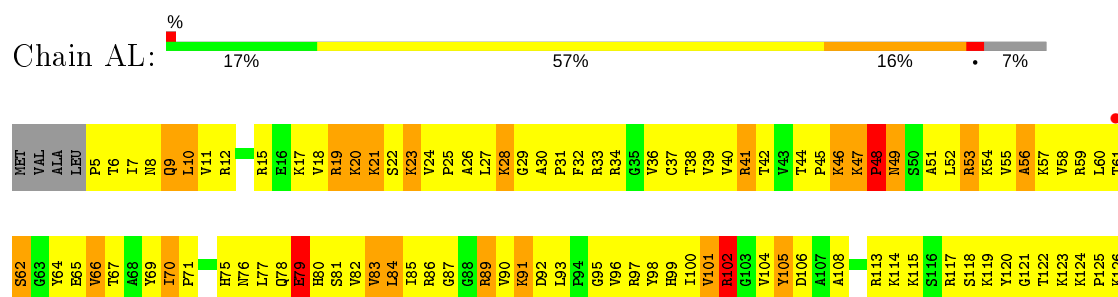
• Molecule 11: 30S RIBOSOMAL PROTEIN S11

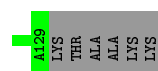


• Molecule 11: 30S RIBOSOMAL PROTEIN S11

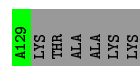
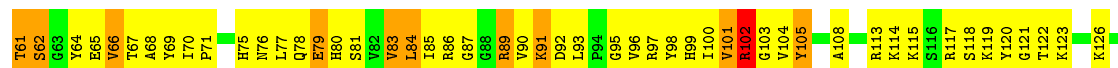


• Molecule 12: 30S RIBOSOMAL PROTEIN S12





• Molecule 12: 30S RIBOSOMAL PROTEIN S12



• Molecule 13: 30S RIBOSOMAL PROTEIN S13



• Molecule 13: 30S RIBOSOMAL PROTEIN S13



• Molecule 14: 30S RIBOSOMAL PROTEIN S14

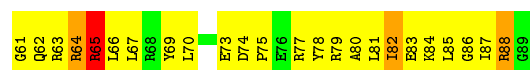




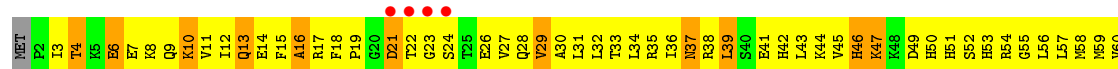
• Molecule 14: 30S RIBOSOMAL PROTEIN S14



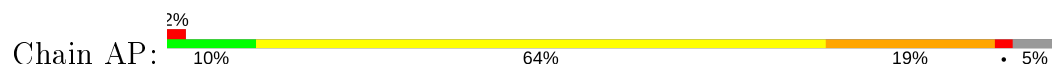
• Molecule 15: 30S RIBOSOMAL PROTEIN S15



• Molecule 15: 30S RIBOSOMAL PROTEIN S15



• Molecule 16: 30S RIBOSOMAL PROTEIN S16



• Molecule 16: 30S RIBOSOMAL PROTEIN S16

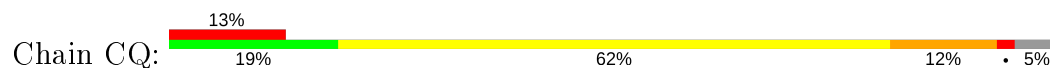




• Molecule 17: 30S RIBOSOMAL PROTEIN S17



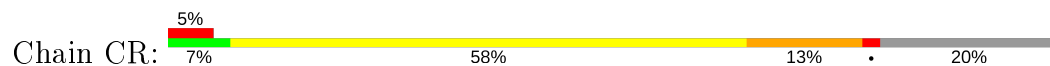
• Molecule 17: 30S RIBOSOMAL PROTEIN S17



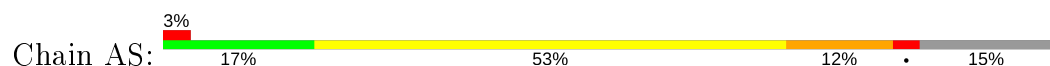
• Molecule 18: 30S RIBOSOMAL PROTEIN S18

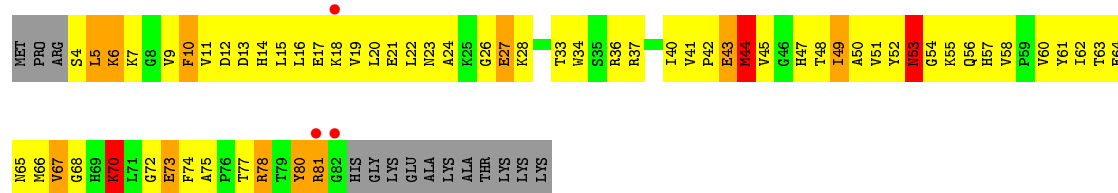


• Molecule 18: 30S RIBOSOMAL PROTEIN S18

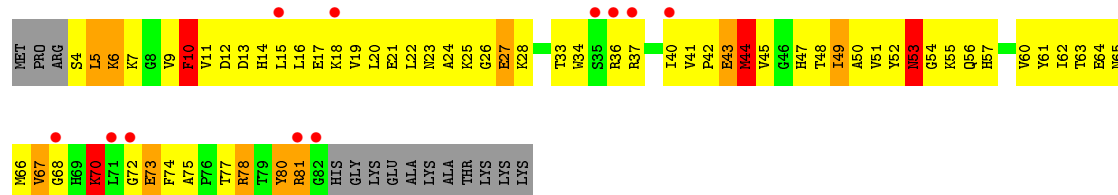
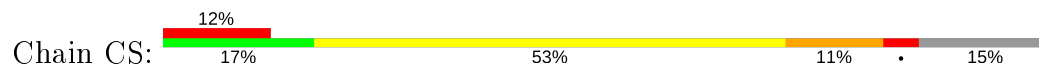


• Molecule 19: 30S RIBOSOMAL PROTEIN S19

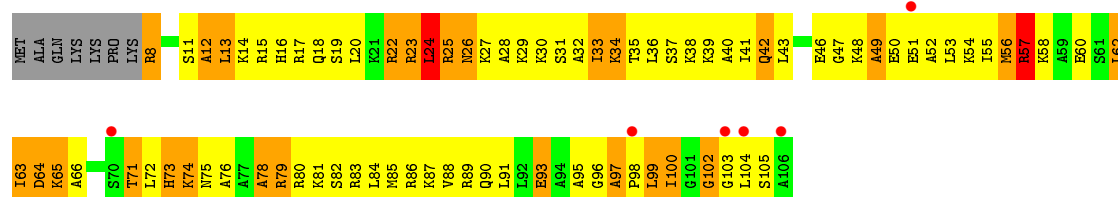
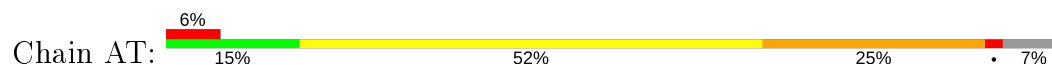




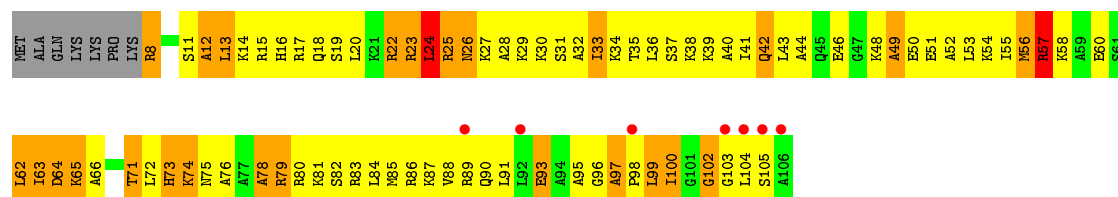
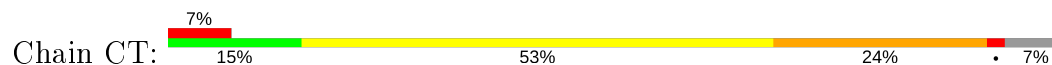
• Molecule 19: 30S RIBOSOMAL PROTEIN S19



• Molecule 20: 30S RIBOSOMAL PROTEIN S20



• Molecule 20: 30S RIBOSOMAL PROTEIN S20



• Molecule 21: 30S RIBOSOMAL PROTEIN THX

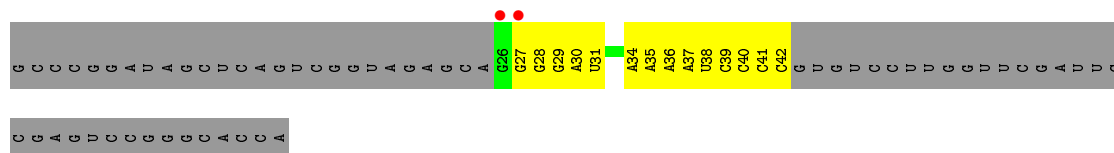


• Molecule 21: 30S RIBOSOMAL PROTEIN THX

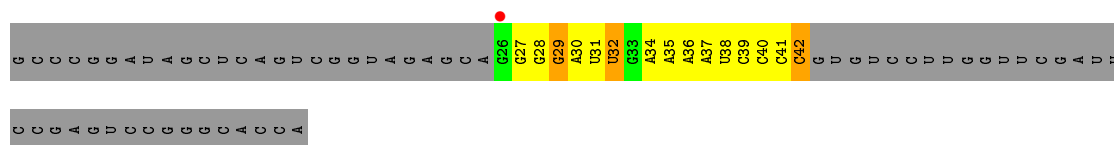




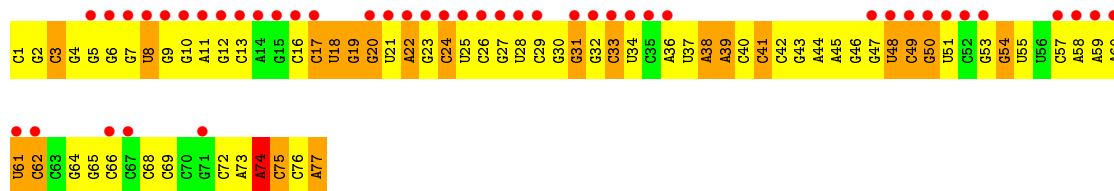
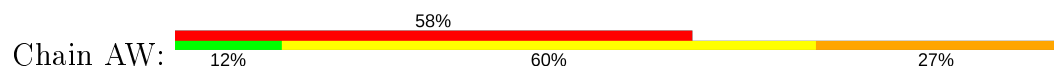
• Molecule 22: P-SITE RNA ASL-PHE



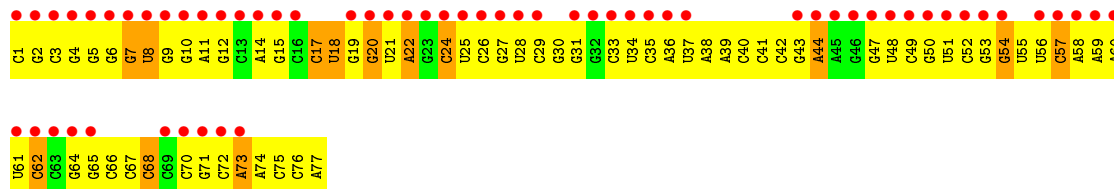
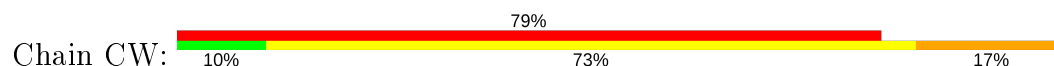
• Molecule 22: P-SITE RNA ASL-PHE



• Molecule 23: TRNA-FMET



• Molecule 23: TRNA-FMET



• Molecule 24: MRNA



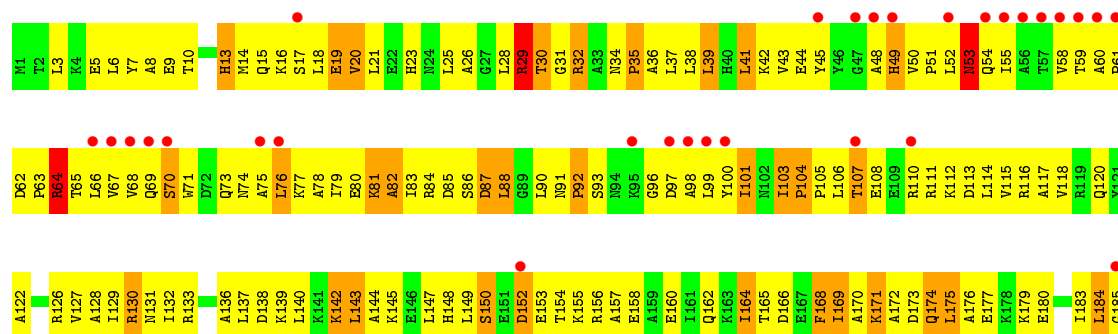
• Molecule 24: MRNA

Chain CX:  29% 6% 65%

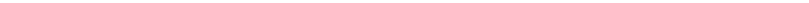


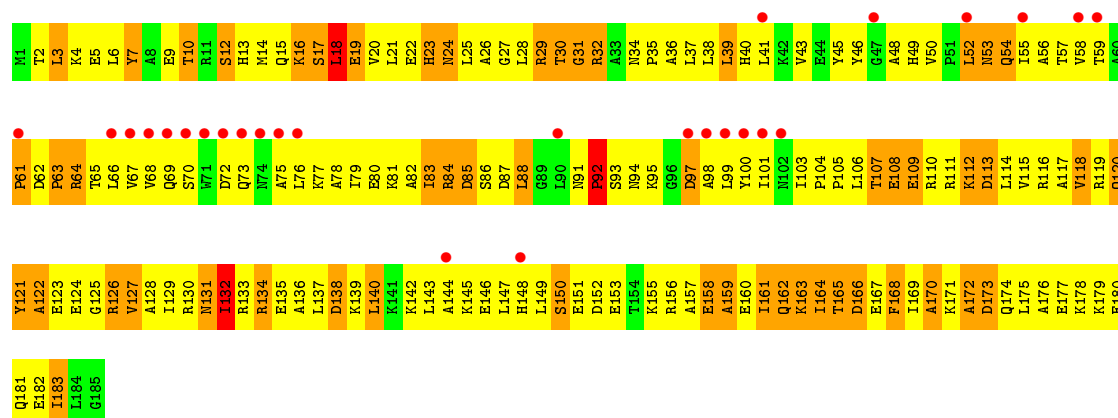
- Molecule 25: RIBOSOME RECYCLING FACTOR

Chain AY: 



- Molecule 25: RIBOSOME RECYCLING FACTOR

Chain CY: 



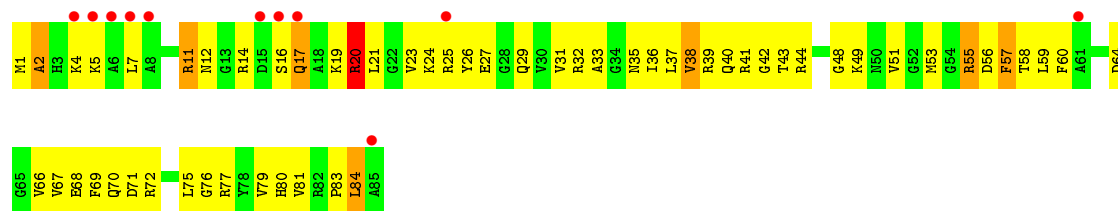
- Molecule 26: 50S RIBOSOMAL PROTEIN L27

Chain B0: 

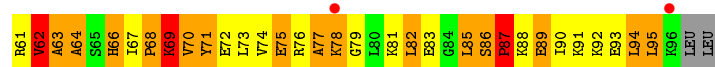


● Molecule 26: 50S RIBOSOMAL PROTEIN L27

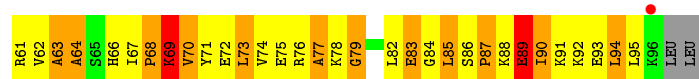
Chain D0: 



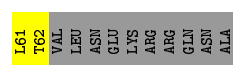
• Molecule 27: 50S RIBOSOMAL PROTEIN L28



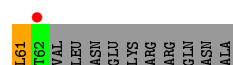
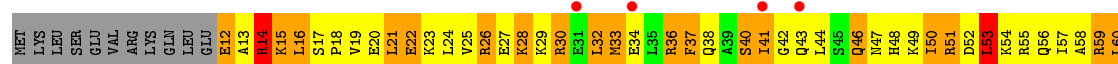
• Molecule 27: 50S RIBOSOMAL PROTEIN L28



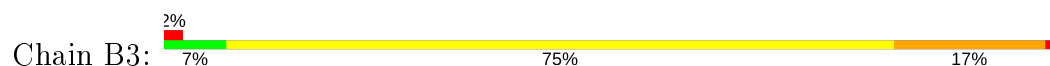
• Molecule 28: 50S RIBOSOMAL PROTEIN L29

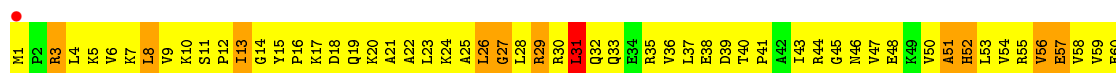


• Molecule 28: 50S RIBOSOMAL PROTEIN L29

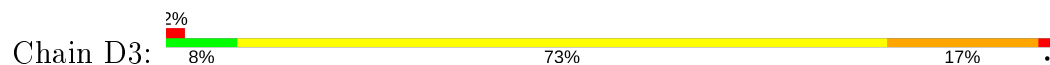


• Molecule 29: 50S RIBOSOMAL PROTEIN L30





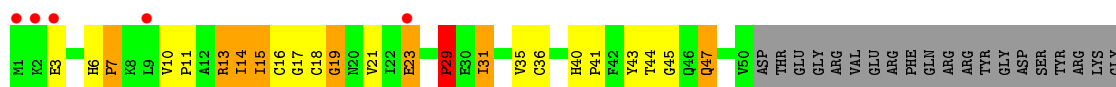
• Molecule 29: 50S RIBOSOMAL PROTEIN L30



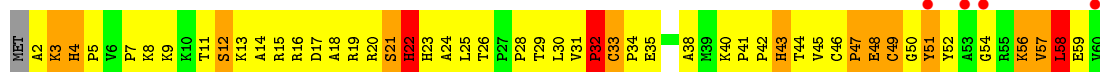
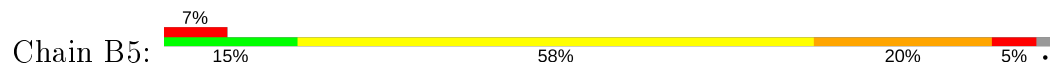
• Molecule 30: 50S RIBOSOMAL PROTEIN L31



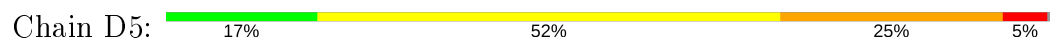
• Molecule 30: 50S RIBOSOMAL PROTEIN L31



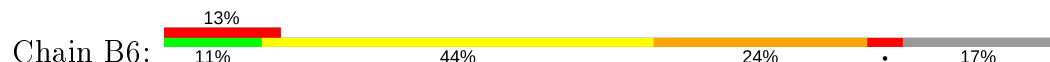
• Molecule 31: 50S RIBOSOMAL PROTEIN L32



• Molecule 31: 50S RIBOSOMAL PROTEIN L32



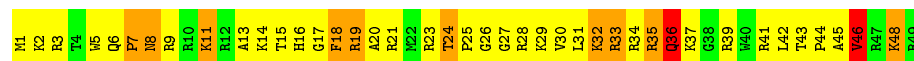
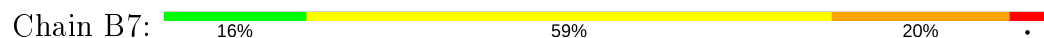
• Molecule 32: 50S RIBOSOMAL PROTEIN L33



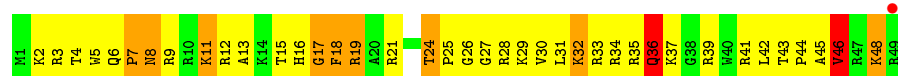
• Molecule 32: 50S RIBOSOMAL PROTEIN L33



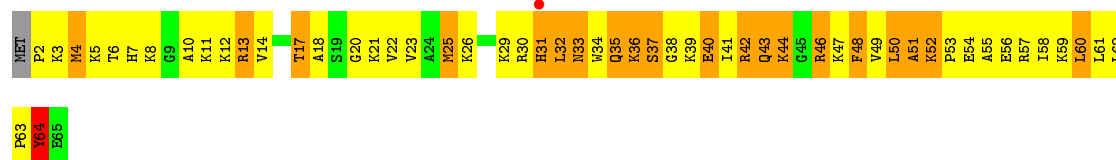
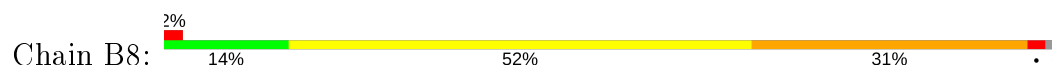
- Molecule 33: 50S RIBOSOMAL PROTEIN L34



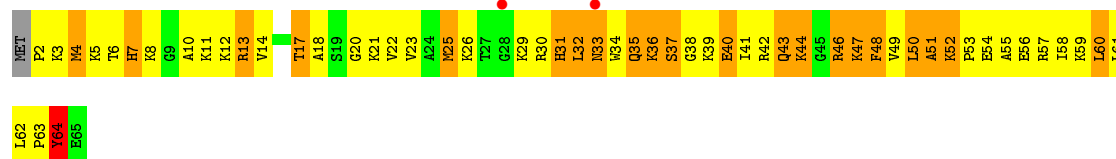
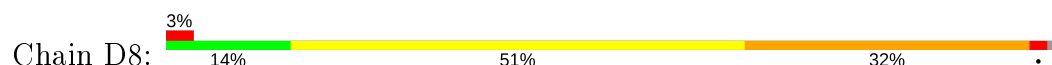
- Molecule 33: 50S RIBOSOMAL PROTEIN L34



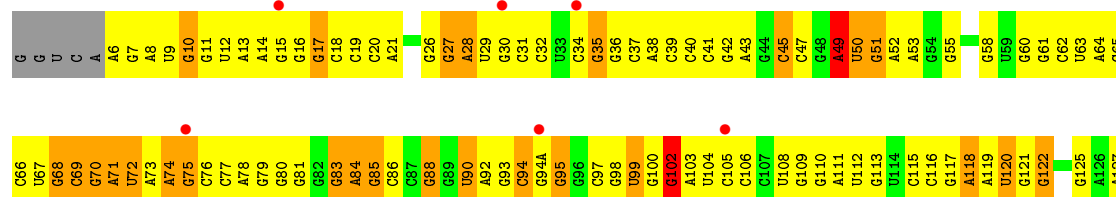
- Molecule 34: 50S RIBOSOMAL PROTEIN L35



- Molecule 34: 50S RIBOSOMAL PROTEIN L35



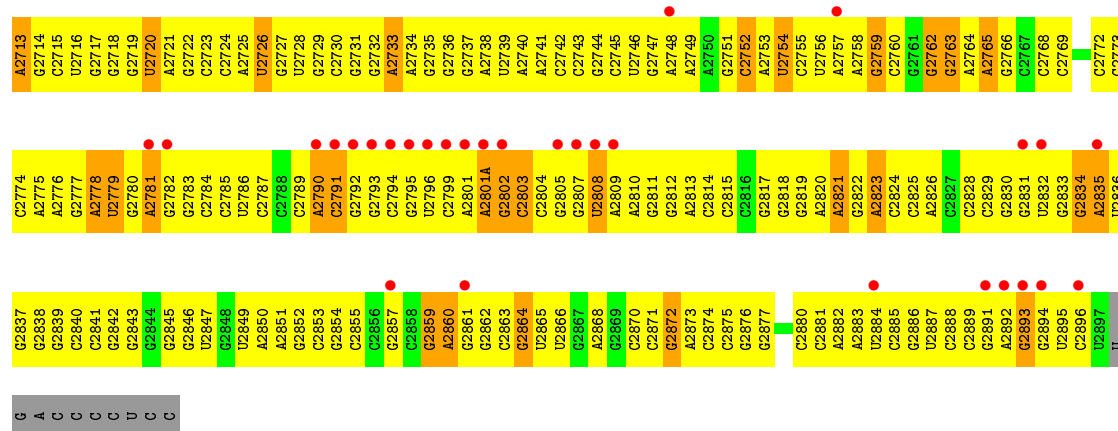
- Molecule 35: 23S ribosomal RNA



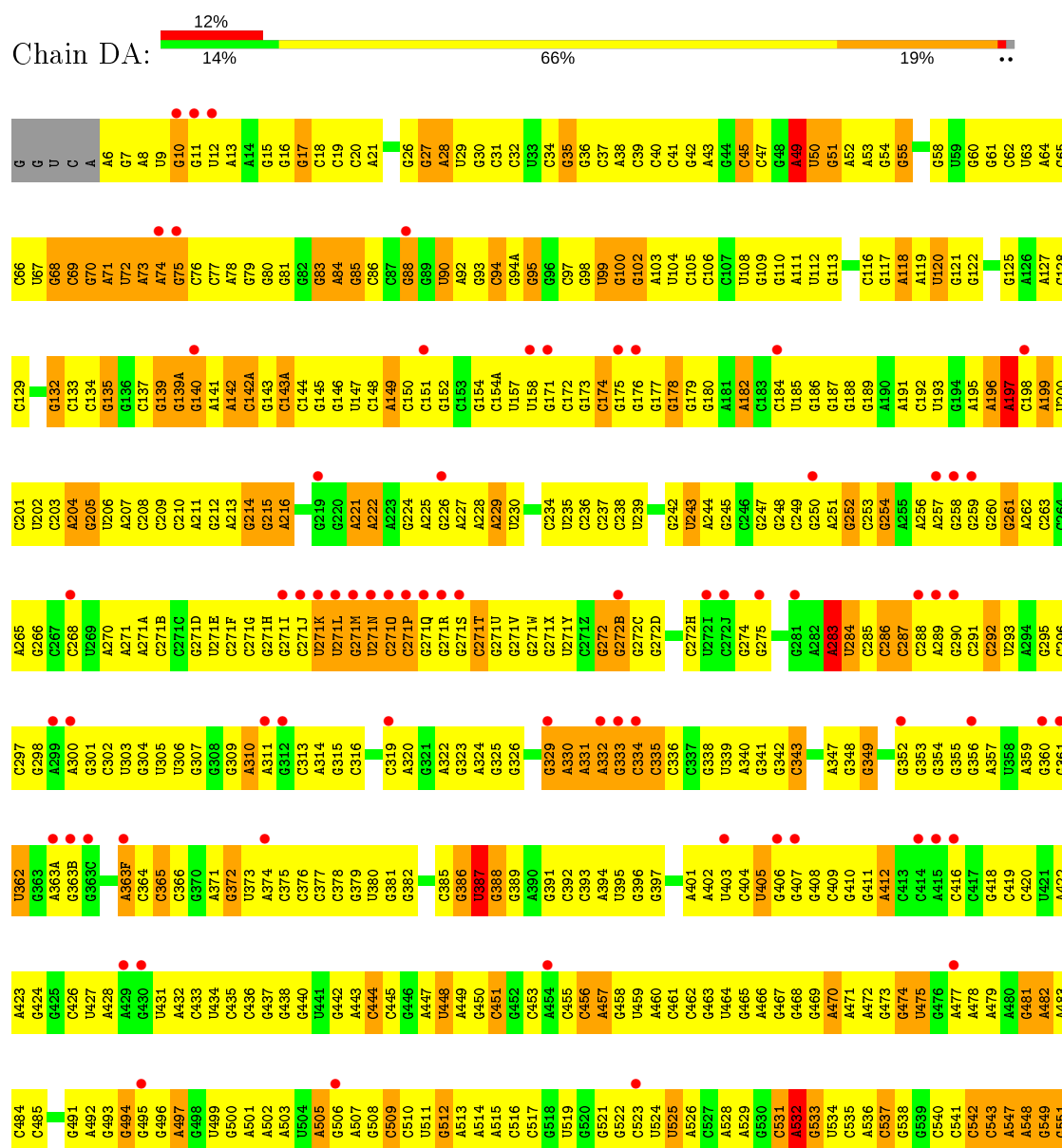


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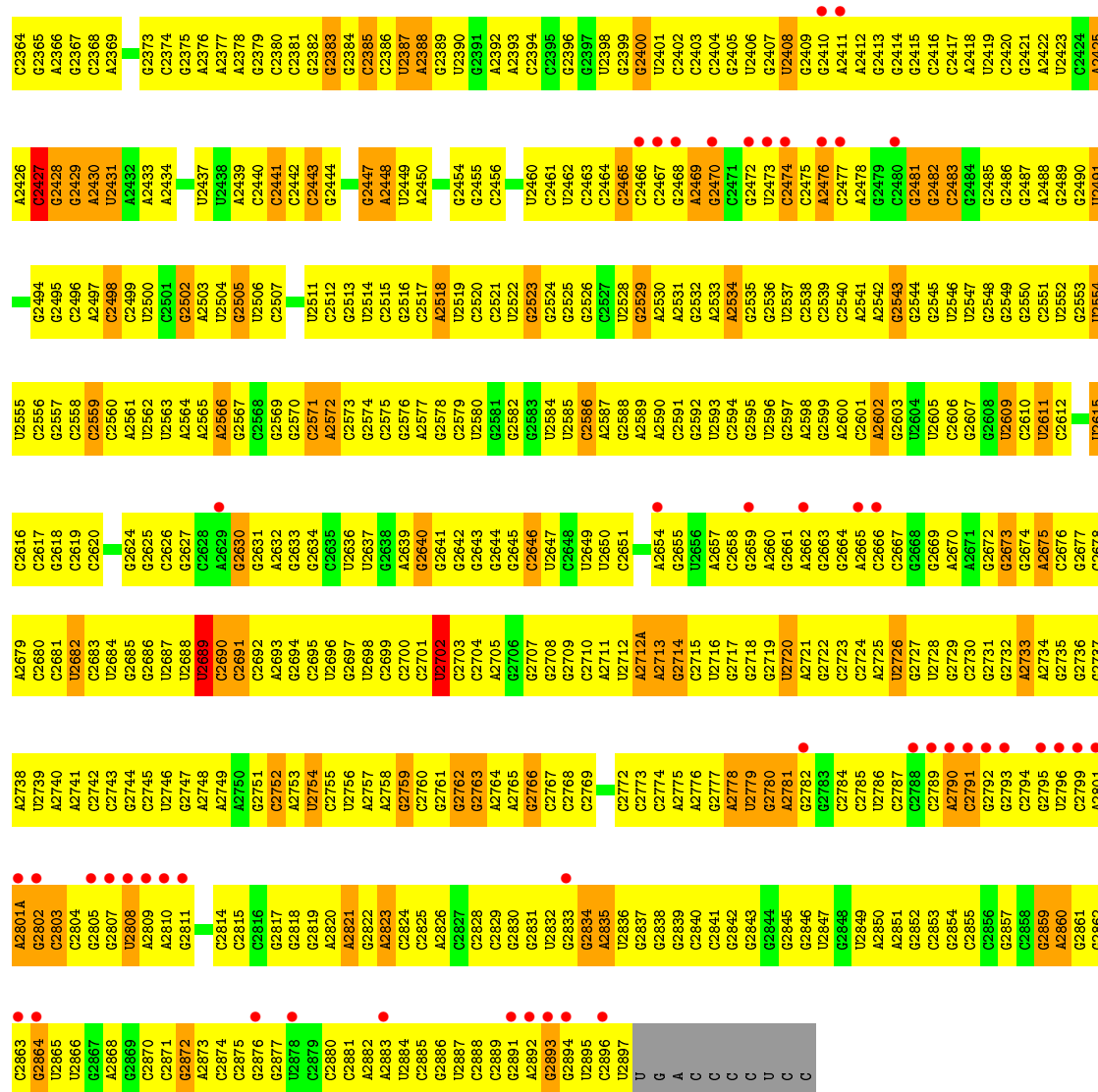


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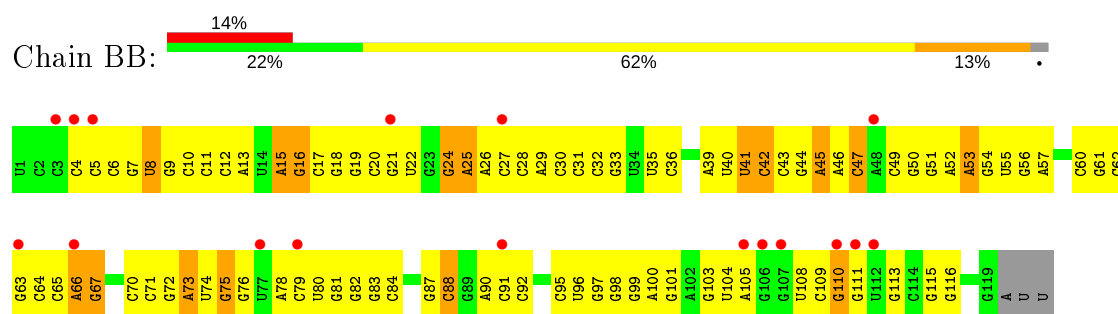


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G1418	A1353	C1291	C1230	G1169	C1052	C991	G932	A863	A802	C679	G614B	U555
A1419	A1354	U1292	G1231	G1170	A1106	G992	A933	U803	U804	G680	G615	U556
U1420	G1355	C1293	G1232	G1171	G1107	G993	G934	C965	G805	G681	G616	U557
G1421	G1356		C1233	G1172	G1108	C994	C935	A866	G806	G682	G617	U558
G1422	U1357	C1297	U1234	U1173	C1109	C995	C936	C967	U807	C683	G618	U559
G1423	G1358	C1298		U1174	G1110	A996	U937	U868	G808	G684	G619	U560
A1424	A1359	U1299	A1237	G1175	G1111	G997	G938	A869	G809	G685	G620	U561
G1425	A1360	U1300	G1238	A1176	A1112	C998	G939	A870	G810	G686	G621	U562
G1426	G1362	A1301	G1239	A1177	G1113	U999	G940	U871	U810	G687	G622	U563
A1427	C1363	G1302	U1240	C1178	U1114	G1002	A941	A872	U811	U688	G623	U564
G1428	G1364	G1303	C1241	C1179	G1115	G1003	G942	G873	C812	U689	G624	U565
G1429	A1365	C1304	A1242	C1180	G1116	C1004	U943	G874	U813	G690	G625	U566
C1430	G1366	C1305	G1243	C1181	G1117	G1005	G944	C875	C814	C691	U626	U568
U1431	A1367	C1306	G1244	G1183	G1118	C1006	A945	U876	C815	G692	G627	U569
C1432	A1367	A1307	G1245	G1184		C1007	G946	C877	C816	C693	G628	U570
U1433	G1368	A1308	A1246	C1185	G1122	C1008	G947	A878	C817	U694	G629	A571
U1434	G1369	G1309	A1247	G1186	G1123	A1009	G948	G879	U818	G695	G630	A572
G1435	C1370	G1310	G1248	G1187	C1124	G1010	C949	G880	A819	G696	A631	U573
U1436	G1371	G1311	U1249	U1188	G1127	G1011	G950	G881	A820	C697	A632	C574
G1437	U1372	U1312	G1250	A1189		U1012	C951	G882	A821		A633	A575
U1438	A1373	C1313	C1251	G1190	A1127	G1013	G952	G883	U822	G700	C634	U576
A1439	G1374	G1314	G1252	G1191	A1128	C1014	A953	C884	G823	G701	C635	U577
G1440	C1375	C1315	A1253	G1192	A1129	U1015	G954	G892	A824	U702	G636	A578
U1441	C1376	A1316	U1254	G1193	U1130	G1016	C955	C893	G825	G703	G637	U579
G1442	G1377	A1317	A1255	A1194	G1131	G1017	G956	C894	U826	U704		C580
A1443	A1378	C1318	G1256	G1195	A1132	U1018	A957	U895	U827	G705	G641	C581
G1444	C1379	G1319	C1257	C1196	U1133	C1018	U958	A896	U828	A706	G642	G582
U1445	G1380	A1320	G1258	G1197	C1135	U1019	A959	C897	A829	G707	A643	G583
A1446	C1381	A1321	G1259	U1198	G1136	A1020	A960	C898	G830	C708	A644	G584
G1447	G1382	A1322	G1260	U1199	G1137	A1021	C961	A899	G831	U709	C645	G585
U1448	C1383	G1323	C1261	C1200	G1138	G1022	G962	A900	U832	G710	A646	A586
G1449	A1384	G1324	A1262	C1201		U1023	C963	C902	U833	G711	G647	C587
U1450	G1385	G1325	U1263	G1202	U1141	G1024	C964	C903	C834	G712	G648	U588
C1450A	U1326	A1326	G1264	G1203	G1142	G1025	C965	C904	A835	G713	G649	C589
C1451	C1327	A1265	A1205	A1204	A1142A	U1026	G966		G836	U714	C650	A590
U1452	G1328	G1266	U1205	A1207	A1143	A1027	C967	U905	C837	G715	G651	C591
U1453	U1329	U1267		G1144	G1144	A1028	G968	G906	C838	A716	C652	G592
G1454	A1330	A1268	C1208	C1145	C1145	A1029	U969	U907	U839	G717	G656	G593
U1455	G1331	A1269	G1209	C1146		G1030	C970	C908	U840	A718	U657	U594
G1456	C1332	C1270	A1210	G1149	G1149	G1031	C971	A909	A841	C719	C658	C595
A1457	C1333	A1271	U1211	C1150		A1032	G972	A910	U842		C659	G596
G1458	U1334	A1272	G1212	G1151	U1033	U1033	A973	A911	A782	A722	G660	U597
U1459	C1335	G1273	A1213	C1152	G1034	G1034	G974	C912	A783	G723	G661	G598
A1460	A1336	A1274	A1214	C1153	U1035	U1035	C975	U913	A784	U724	G662	G599
G1461	G1337	A1275	G1215	G1154	G1036	G1036	G975A	C914	G785	G725	G663	G600
C1462	C1338	A1276	G1216	A1155	G1037	G1037	C976	C915	C786	G726	C664	C601
G1463	U1339	G1277	C1217	A1156	C1038	U1038	G977	G916	U787	A727	C665	G602
U1464	G1340	A1278	C1218	G1157	G1039	G1039	G978	A917	A788	G728	G666	A603
G1465	U1341	G1279	G1219	C1158	C1040	C1040	G979	A918	U851	G729	U667	G604
U1466	A1342	G1280	A1220	C1159	G1041	A980	U980	G919	C790	C730	G668	C605
C1467	G1343	C1281	C1221	U1159	G1042	A981	A981	G920	C791	G731	G669	U606
U1468	G1344	U1282	C1221A	G1160	C1043	G982	C982	U921	G792	C732	G670	U607
A1469	C1345	G1283	G1222	C1161	G1044	A983	A983	U922	A793	G733	C671	A608
G1470	G1346	A1284	G1223	G1162	A1045	A984	A984	C923	G794		C672	A609
A1471	C1347	G1285	C1224	G1163	A1046	C985	C985	C924	C795	C736	G673	G610
U1472	G1348	A1286	G1225	G1164	A1047	C986	C986	U858	G796		G674	C611
	A1349	A1287	A1226	U1165		A1048	G987	G859	C797			

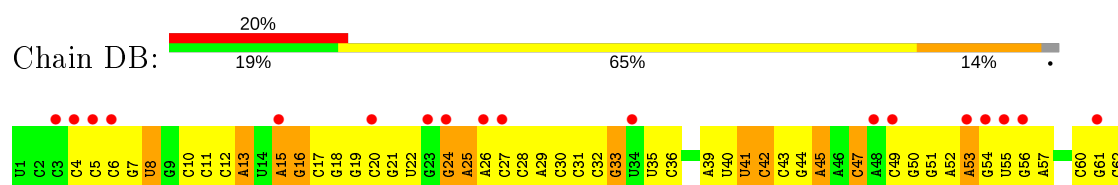


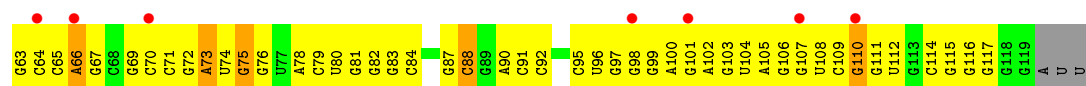


• Molecule 36: 5S ribosomal RNA

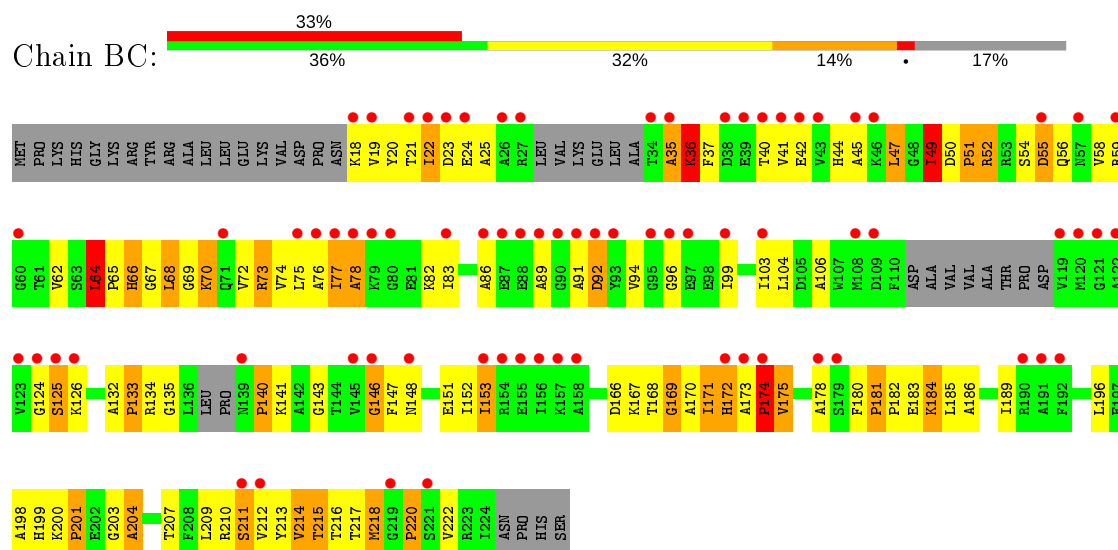


• Molecule 36: 5S ribosomal RNA

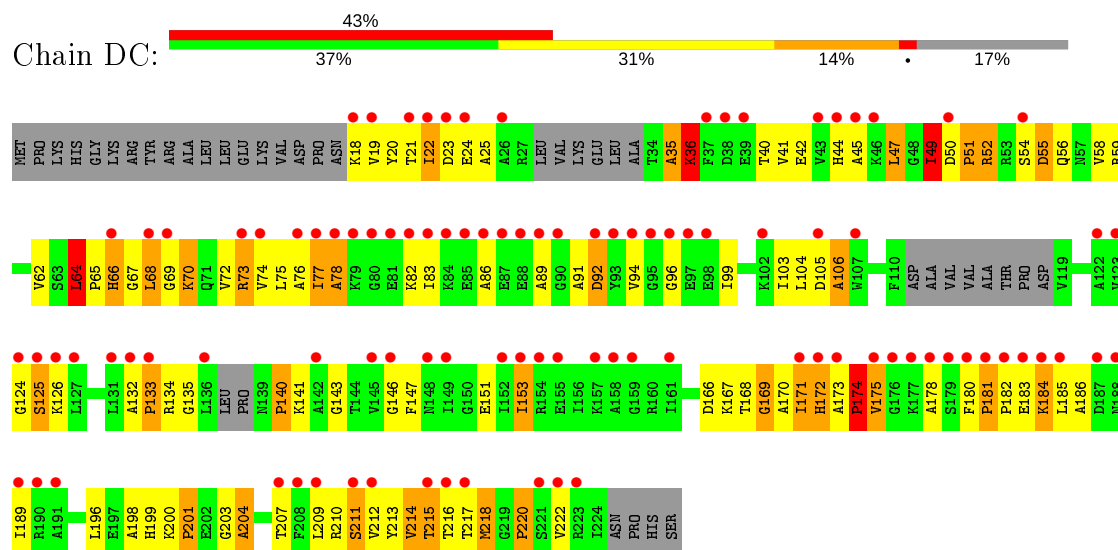




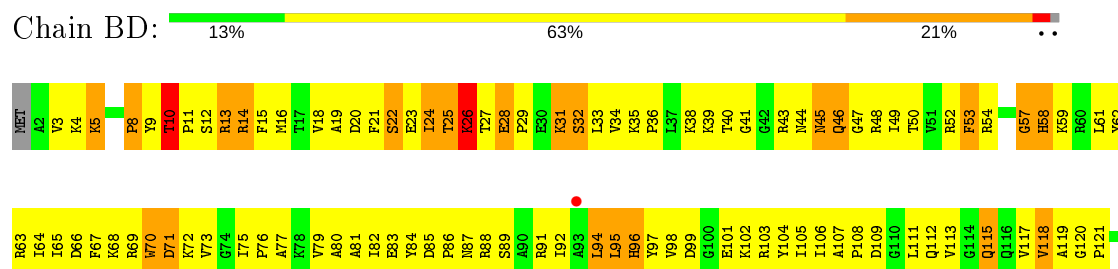
• Molecule 37: 50S RIBOSOMAL PROTEIN L1

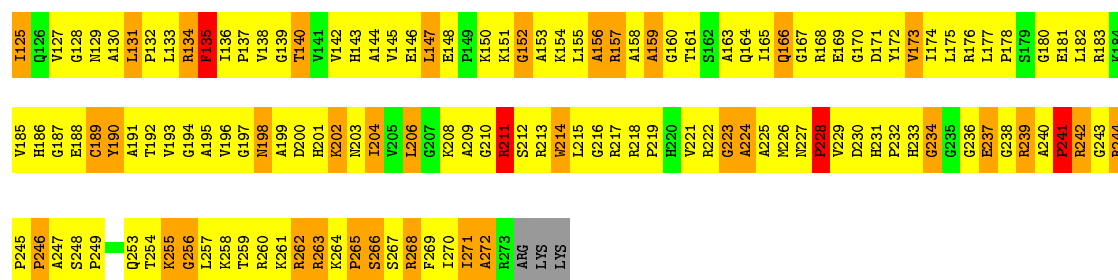


• Molecule 37: 50S RIBOSOMAL PROTEIN L1

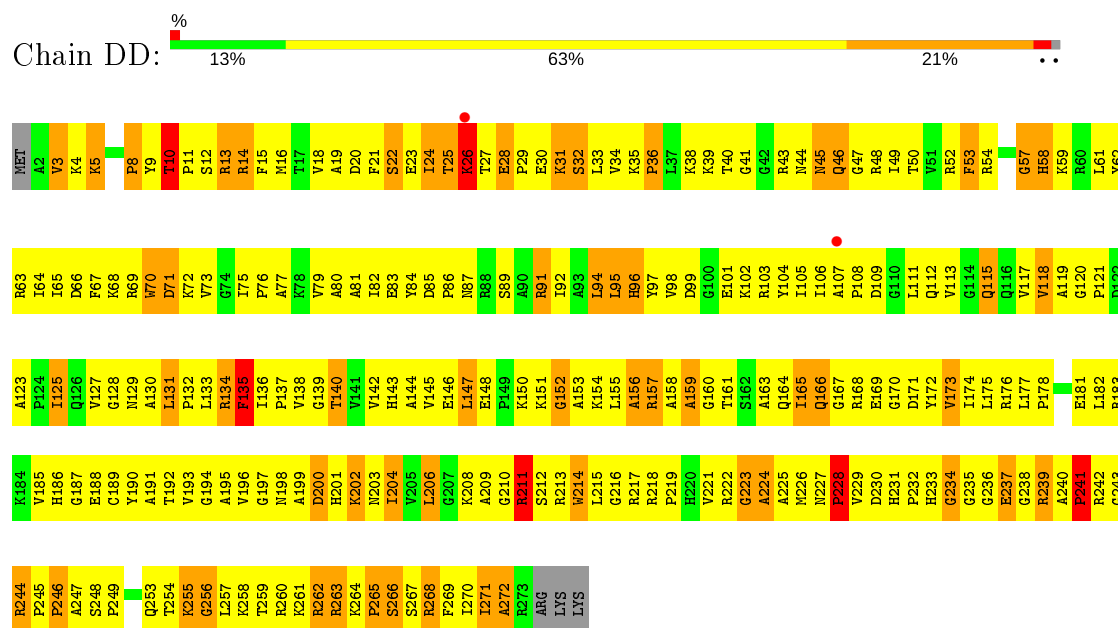


• Molecule 38: 50S RIBOSOMAL PROTEIN L2

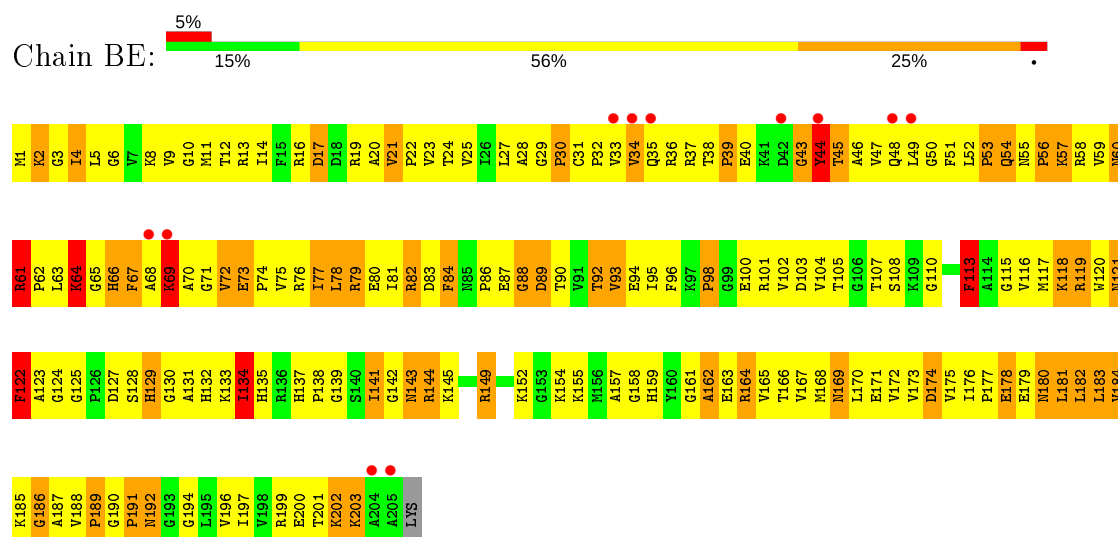




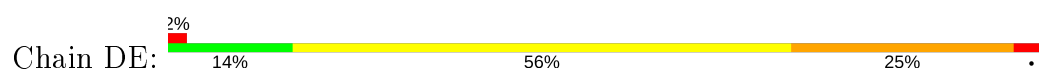
• Molecule 38: 50S RIBOSOMAL PROTEIN L2

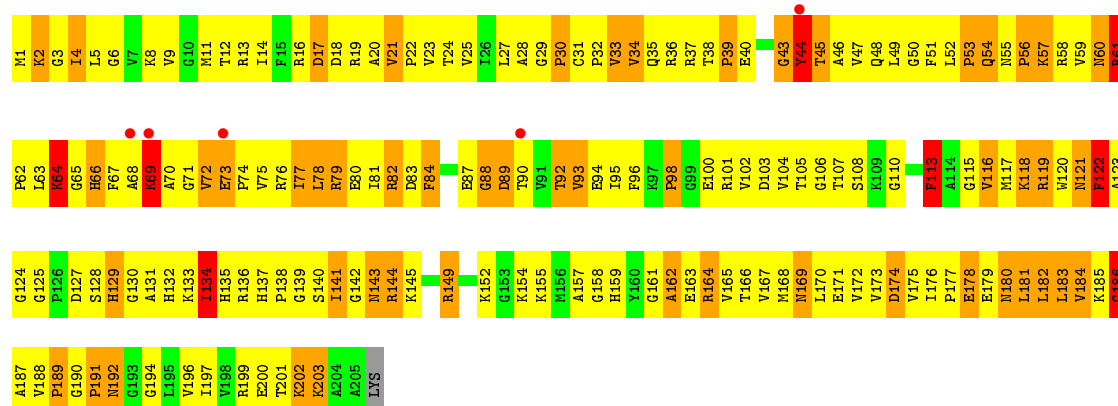


• Molecule 39: 50S RIBOSOMAL PROTEIN L3

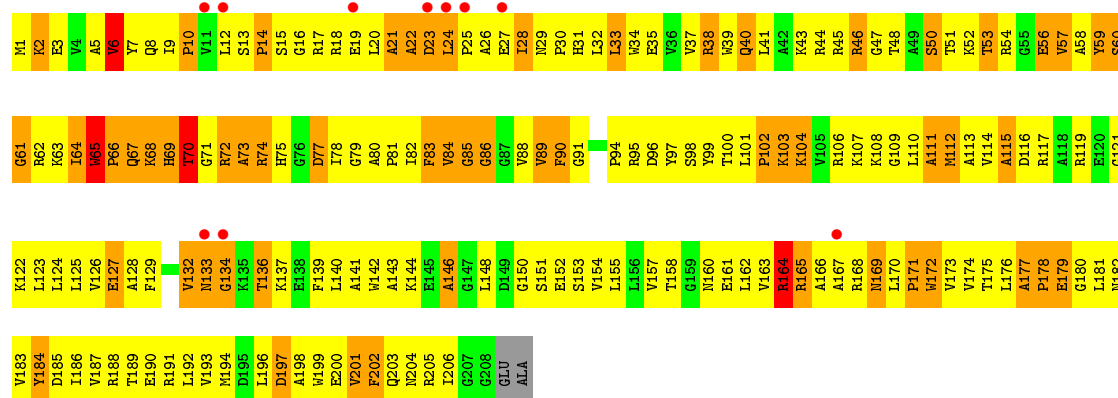
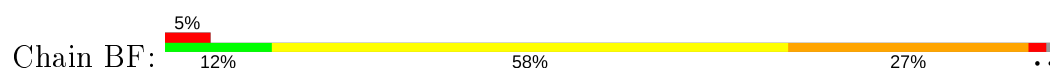


• Molecule 39: 50S RIBOSOMAL PROTEIN L3

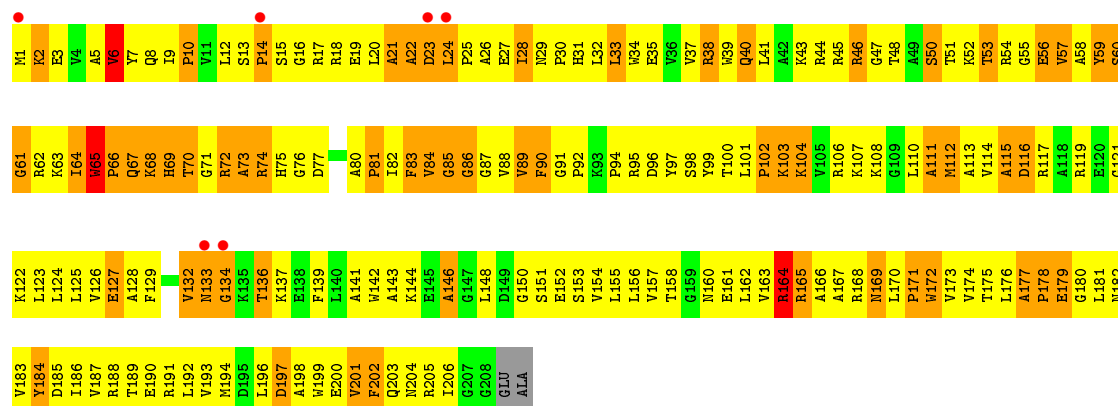
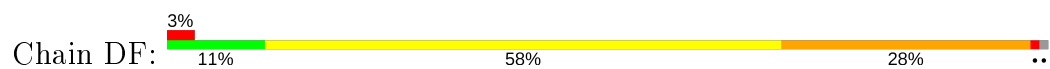




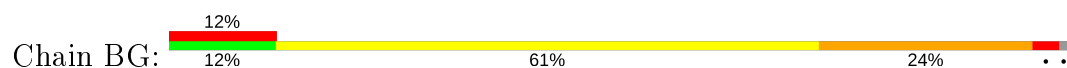
• Molecule 40: 50S RIBOSOMAL PROTEIN L4

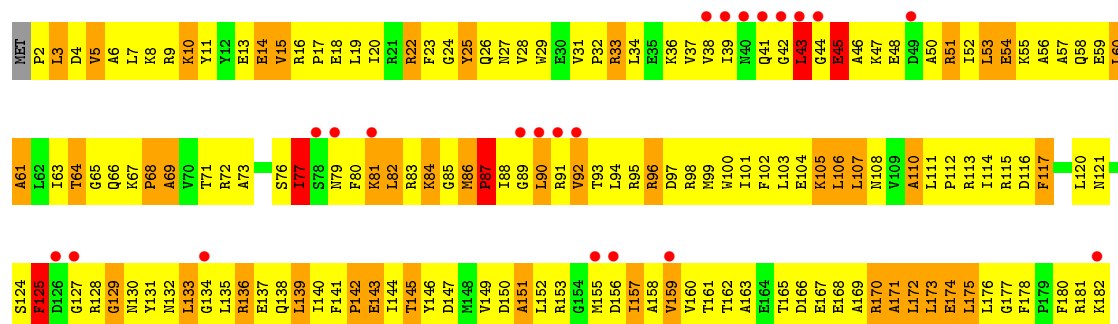


• Molecule 40: 50S RIBOSOMAL PROTEIN L4



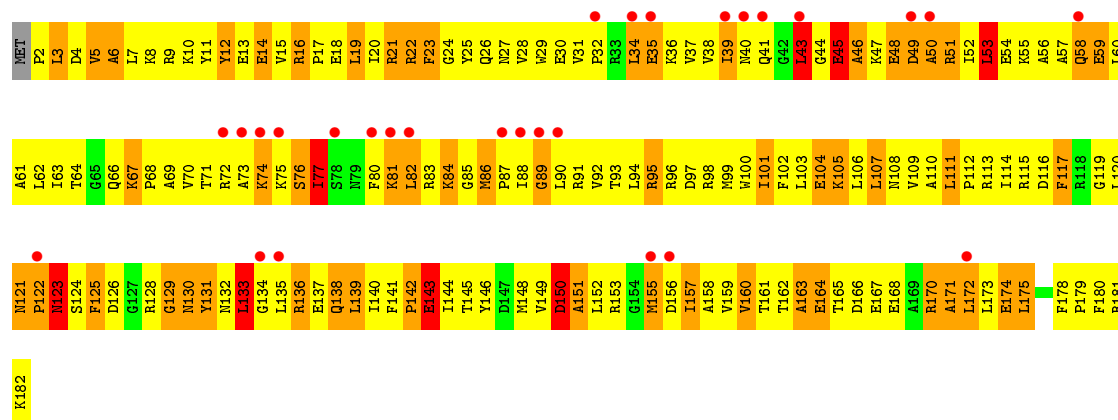
• Molecule 41: 50S RIBOSOMAL PROTEIN L5





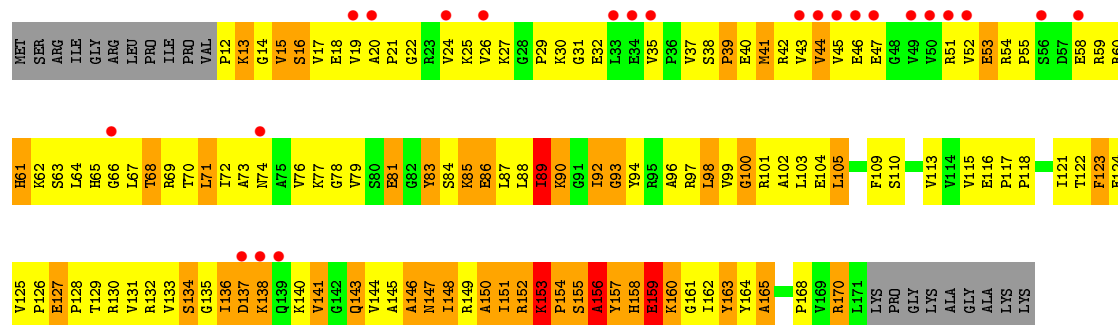
• Molecule 41: 50S RIBOSOMAL PROTEIN L5

Chain DG: 7% 15% 58% 31%



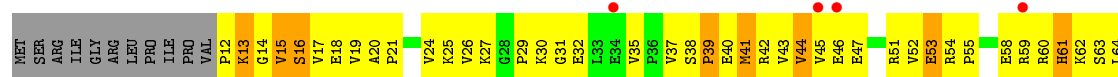
• Molecule 42: 50S RIBOSOMAL PROTEIN L6

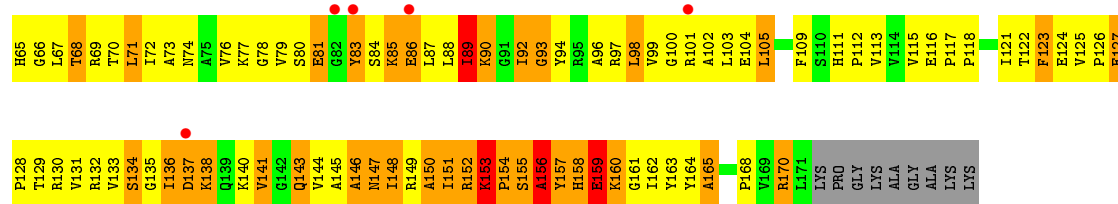
Chain BH: 13% 16% 47% 23% 11%



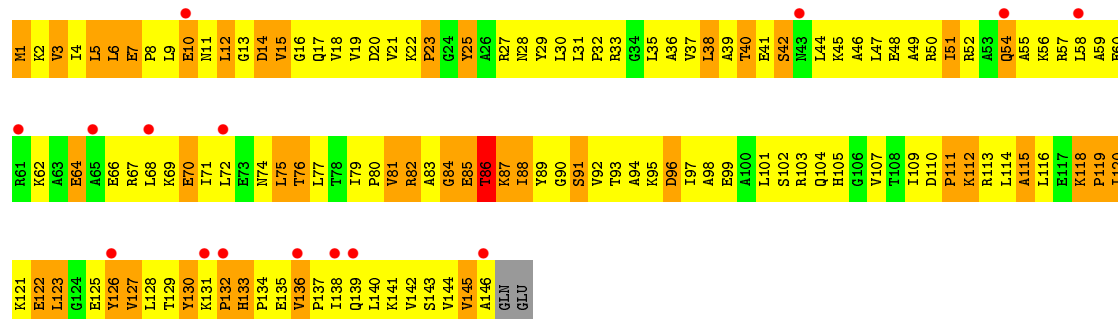
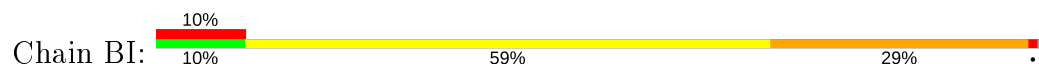
• Molecule 42: 50S RIBOSOMAL PROTEIN L6

Chain DH: 5% 16% 49% 22% 11%

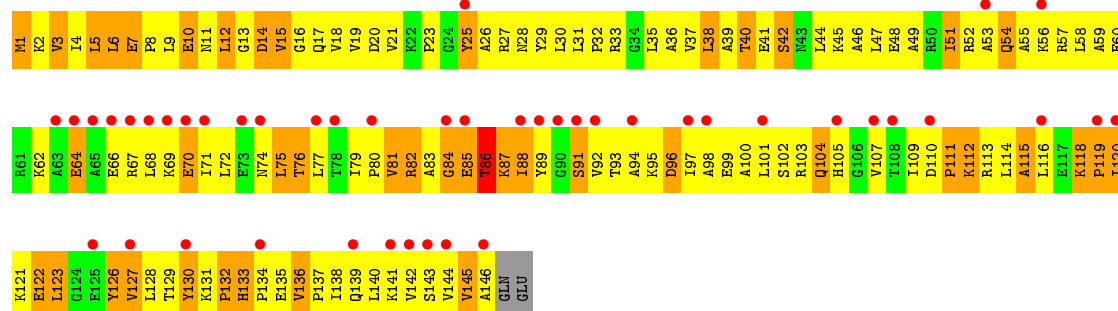
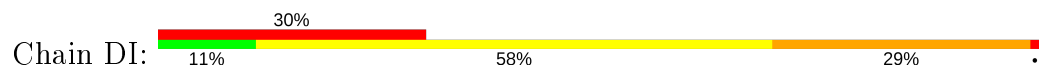




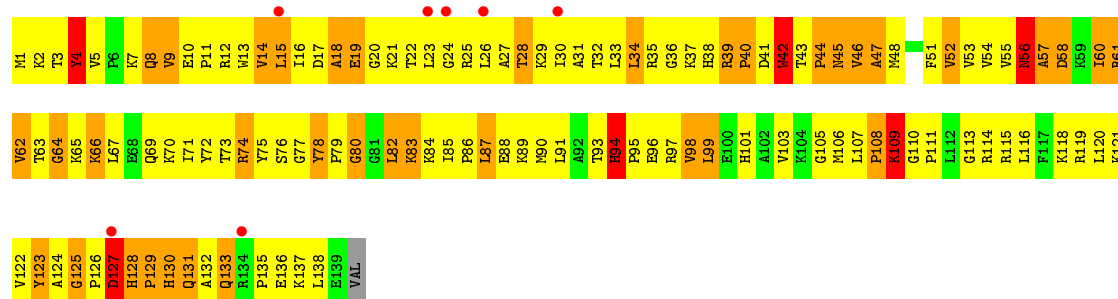
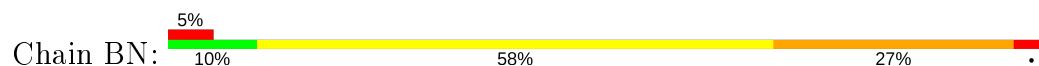
● Molecule 43: 50S RIBOSOMAL PROTEIN L9



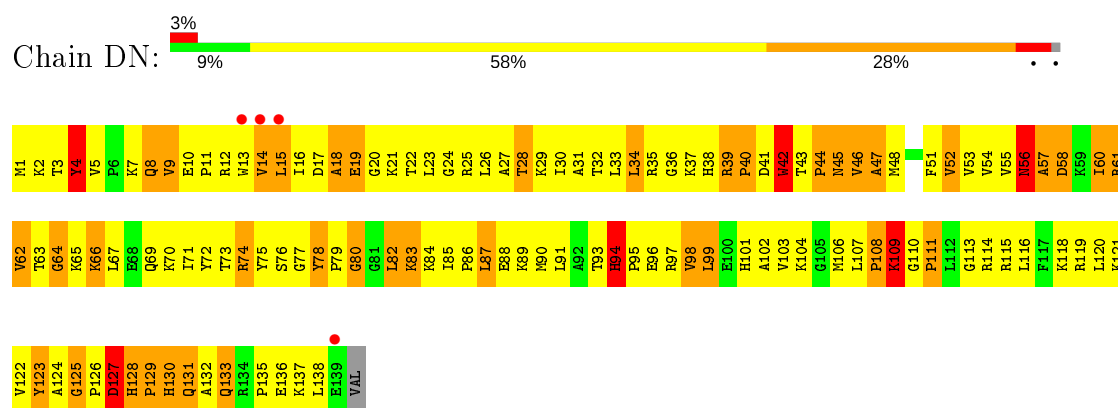
● Molecule 43: 50S RIBOSOMAL PROTEIN L9



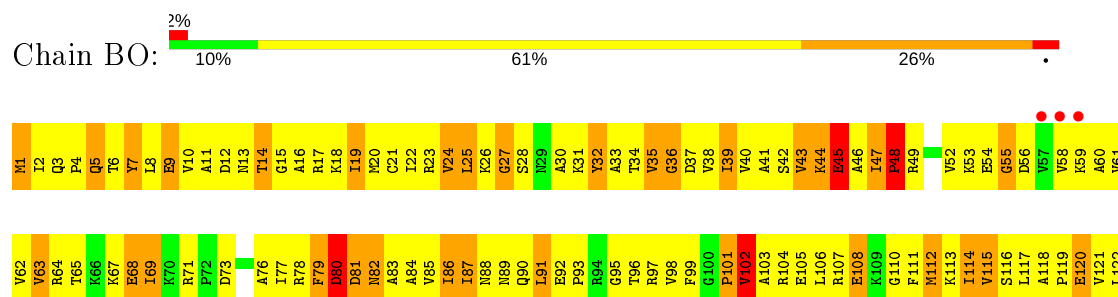
● Molecule 44: 50S RIBOSOMAL PROTEIN L13



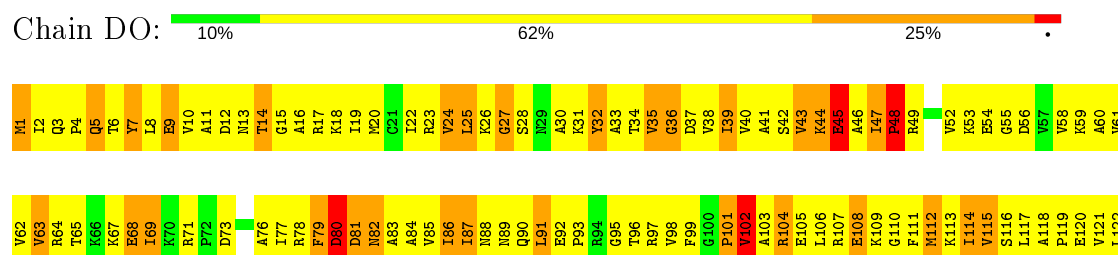
● Molecule 44: 50S RIBOSOMAL PROTEIN L13



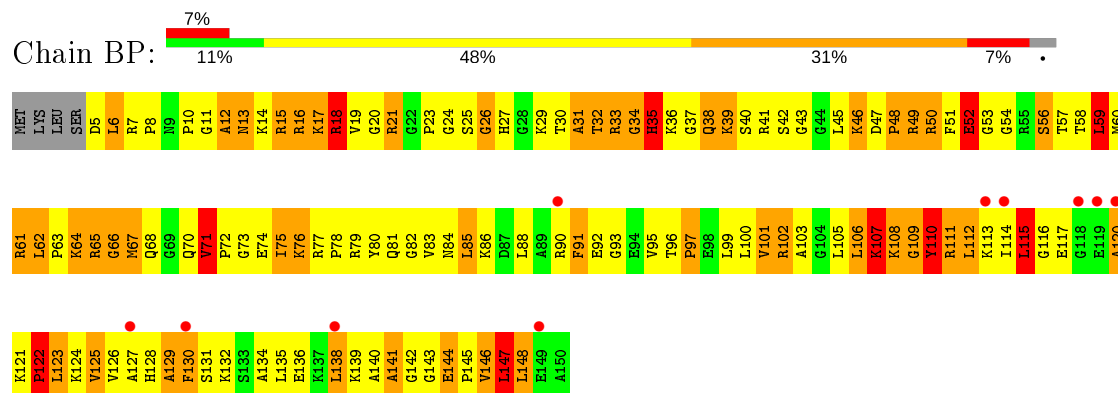
• Molecule 45: 50S RIBOSOMAL PROTEIN L14



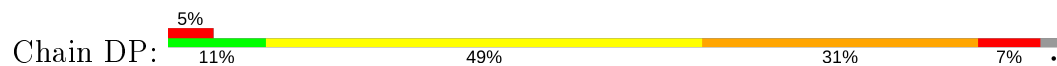
• Molecule 45: 50S RIBOSOMAL PROTEIN L14

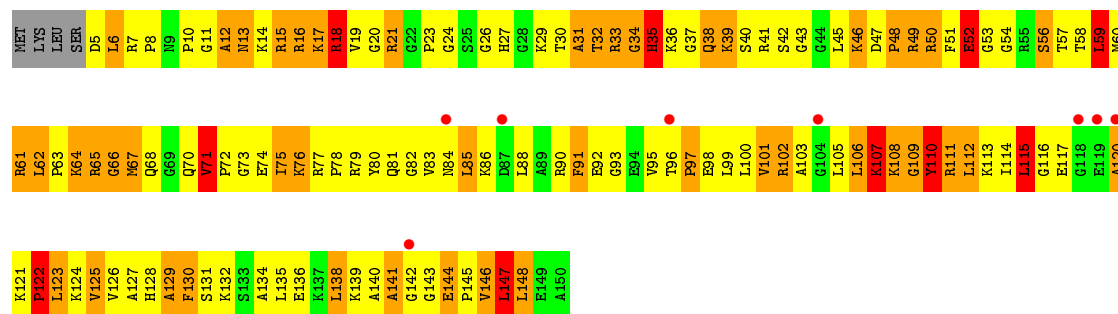


• Molecule 46: 50S RIBOSOMAL PROTEIN L15

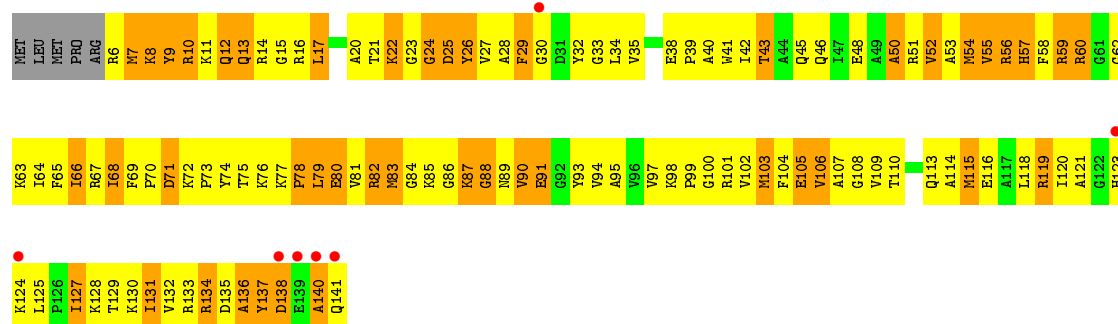
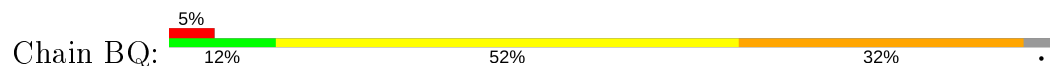


• Molecule 46: 50S RIBOSOMAL PROTEIN L15

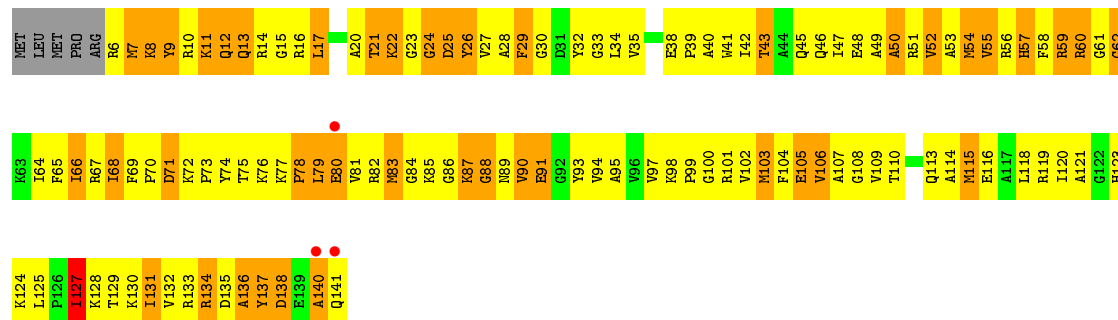
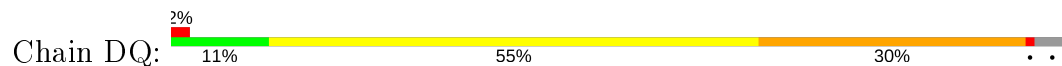




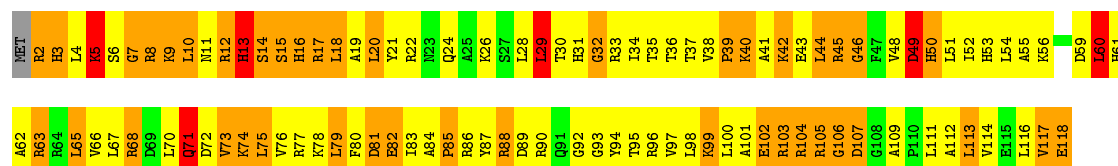
• Molecule 47: 50S RIBOSOMAL PROTEIN L16



• Molecule 47: 50S RIBOSOMAL PROTEIN L16

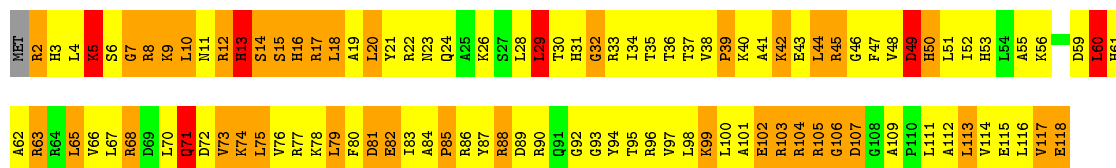


• Molecule 48: 50S RIBOSOMAL PROTEIN L17



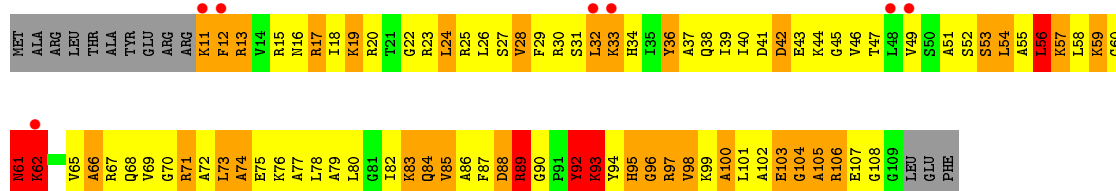
• Molecule 48: 50S RIBOSOMAL PROTEIN L17

Chain DR: 



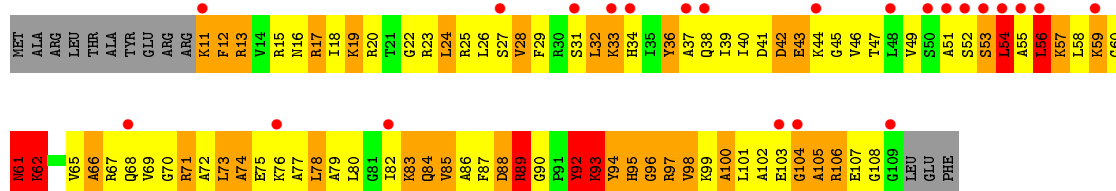
• Molecule 49: 50S RIBOSOMAL PROTEIN L18

Chain BS: 




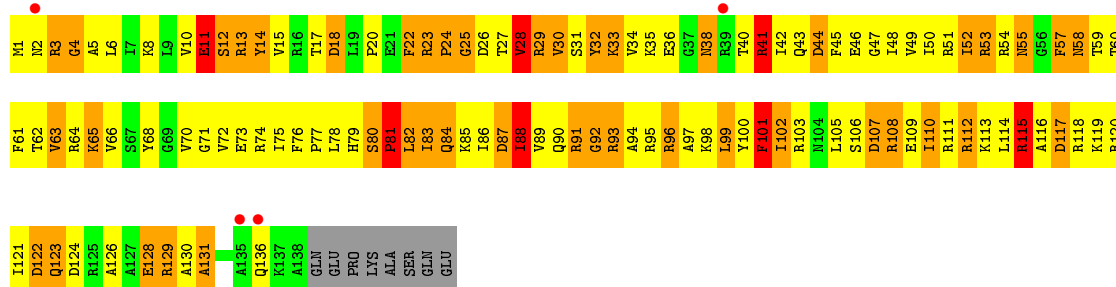
• Molecule 49: 50S RIBOSOMAL PROTEIN L18

Chain DS: 




• Molecule 50: 50S RIBOSOMAL PROTEIN L19

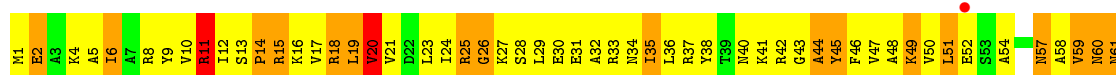
Chain BT: 



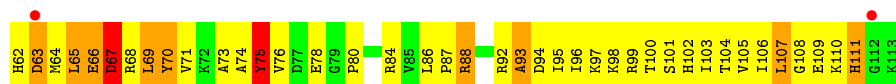
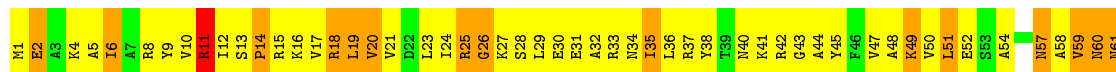
• Molecule 50: 50S RIBOSOMAL PROTEIN L19

Chain DT: 





• Molecule 53: 50S RIBOSOMAL PROTEIN L22



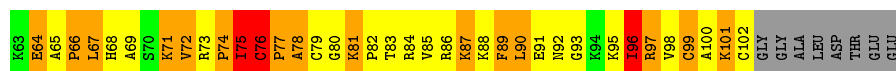
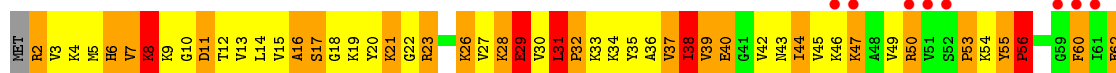
• Molecule 54: 50S RIBOSOMAL PROTEIN L23



• Molecule 54: 50S RIBOSOMAL PROTEIN L23

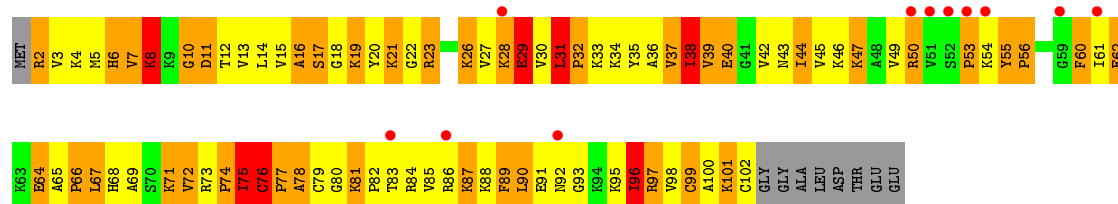


• Molecule 55: 50S RIBOSOMAL PROTEIN L24

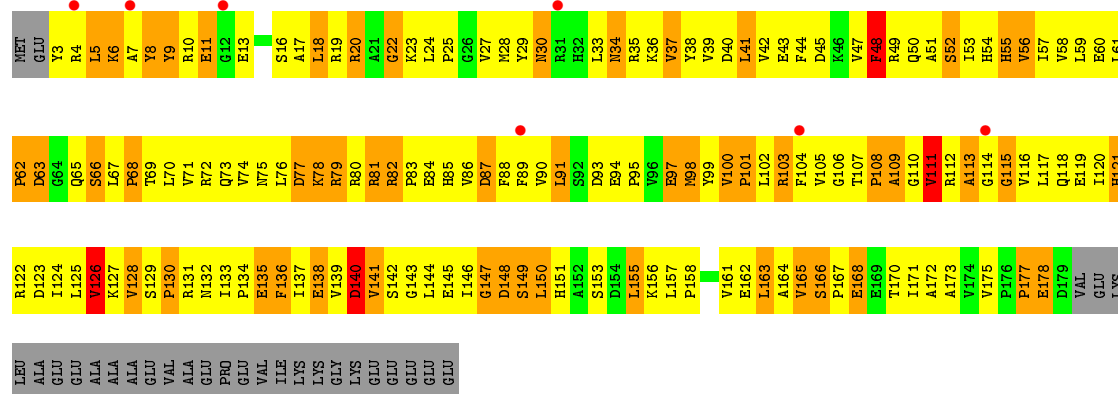


• Molecule 55: 50S RIBOSOMAL PROTEIN L24

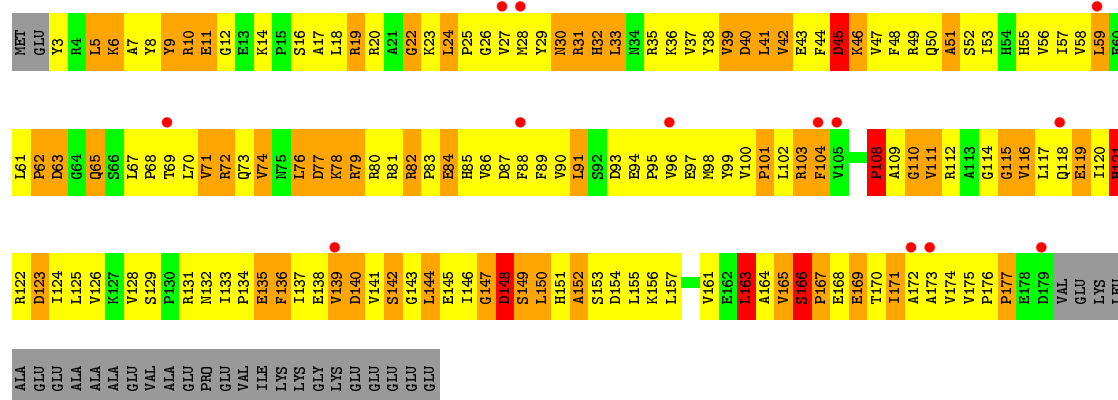




● Molecule 56: 50S RIBOSOMAL PROTEIN L25



● Molecule 56: 50S RIBOSOMAL PROTEIN L25



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	212.41Å 450.11Å 630.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 49.89 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.9 (20.00-3.50) 98.9 (49.89-3.50)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 3.48Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.265 , 0.326 0.260 , 0.318	Depositor DCC
R_{free} test set	34086 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	81.4	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 94.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.33$, $\langle L^2 \rangle = 0.16$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	290487	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 5MU, ZN, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	AA	0.49	0/36190	0.72	11/56486 (0.0%)
1	CA	0.49	0/36190	0.72	5/56486 (0.0%)
2	AB	0.37	0/1936	0.65	0/2611
2	CB	0.36	0/1936	0.64	0/2611
3	AC	0.37	0/1637	0.63	0/2207
3	CC	0.35	0/1637	0.63	0/2207
4	AD	0.43	0/1733	0.68	0/2318
4	CD	0.42	0/1733	0.68	0/2318
5	AE	0.41	0/1163	0.68	0/1566
5	CE	0.41	0/1163	0.68	0/1566
6	AF	0.36	0/856	0.67	0/1154
6	CF	0.39	0/856	0.68	0/1154
7	AG	0.35	0/1276	0.63	0/1709
7	CG	0.34	0/1276	0.62	0/1709
8	AH	0.36	0/1136	0.66	0/1527
8	CH	0.37	0/1136	0.67	0/1527
9	AI	0.35	0/1027	0.63	0/1372
9	CI	0.34	0/1027	0.62	0/1372
10	AJ	0.38	0/808	0.66	0/1087
10	CJ	0.37	0/808	0.65	0/1087
11	AK	0.37	0/900	0.66	0/1213
11	CK	0.37	0/900	0.67	0/1213
12	AL	0.44	0/987	0.72	0/1322
12	CL	0.42	0/987	0.72	0/1322
13	AM	0.33	0/994	0.59	0/1322
13	CM	0.32	0/994	0.60	0/1322
14	AN	0.39	0/501	0.68	0/664
14	CN	0.37	0/501	0.67	0/664
15	AO	0.36	0/745	0.64	0/992
15	CO	0.36	0/745	0.63	0/992
16	AP	0.47	0/717	0.74	0/965
16	CP	0.43	0/717	0.73	0/965

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AQ	0.41	0/837	0.71	0/1119
17	CQ	0.40	0/837	0.71	0/1119
18	AR	0.39	0/579	0.72	0/768
18	CR	0.41	0/579	0.73	0/768
19	AS	0.36	0/643	0.61	0/867
19	CS	0.36	0/643	0.60	0/867
20	AT	0.34	0/765	0.69	0/1007
20	CT	0.33	0/765	0.69	0/1007
21	AU	0.44	0/213	0.64	0/279
21	CU	0.42	0/213	0.63	0/279
22	AV	0.66	0/405	0.77	0/630
22	CV	0.61	0/405	0.77	1/630 (0.2%)
23	AW	0.43	0/1810	0.73	0/2821
23	CW	0.46	0/1810	0.72	0/2821
24	AX	0.70	0/256	0.93	0/397
24	CX	0.68	0/256	0.91	0/397
25	AY	0.40	0/1497	0.71	0/2017
25	CY	0.44	0/1497	0.72	0/2017
26	B0	0.39	0/660	0.64	0/882
26	D0	0.40	0/660	0.64	0/882
27	B1	0.57	0/700	1.00	1/931 (0.1%)
27	D1	0.55	0/700	0.96	3/931 (0.3%)
28	B2	0.49	0/423	0.97	2/560 (0.4%)
28	D2	0.52	0/423	0.91	1/560 (0.2%)
29	B3	0.39	0/473	0.68	0/636
29	D3	0.40	0/473	0.68	0/636
30	B4	0.43	0/241	0.80	3/334 (0.9%)
30	D4	0.44	0/241	0.80	4/334 (1.2%)
31	B5	0.37	0/473	0.69	0/639
31	D5	0.38	0/473	0.71	0/639
32	B6	0.45	0/387	0.63	0/517
32	D6	0.42	0/387	0.63	0/517
33	B7	0.53	0/427	0.83	0/563
33	D7	0.53	0/427	0.84	1/563 (0.2%)
34	B8	0.51	0/516	0.89	1/681 (0.1%)
34	D8	0.49	0/516	0.88	1/681 (0.1%)
35	BA	0.57	2/66757 (0.0%)	0.76	19/104221 (0.0%)
35	DA	0.61	3/66757 (0.0%)	0.76	19/104221 (0.0%)
36	BB	0.40	0/2853	0.70	0/4451
36	DB	0.39	0/2853	0.70	0/4451
37	BC	0.36	0/1145	0.68	7/1556 (0.4%)
37	DC	0.36	0/1146	0.68	7/1558 (0.4%)
38	BD	0.47	0/2155	0.84	2/2907 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	DD	0.49	0/2155	0.84	2/2907 (0.1%)
39	BE	0.44	0/1597	0.78	1/2155 (0.0%)
39	DE	0.47	0/1597	0.79	1/2155 (0.0%)
40	BF	0.46	1/1659 (0.1%)	0.72	0/2246
40	DF	0.46	1/1659 (0.1%)	0.72	0/2246
41	BG	0.39	0/1498	0.76	2/2013 (0.1%)
41	DG	0.38	0/1498	0.76	0/2013
42	BH	0.37	0/1246	0.72	1/1684 (0.1%)
42	DH	0.40	0/1246	0.73	1/1684 (0.1%)
43	BI	0.39	0/1147	0.72	0/1553
43	DI	0.40	0/1147	0.71	0/1553
44	BN	0.40	0/1132	0.74	0/1527
44	DN	0.45	0/1132	0.75	0/1527
45	BO	0.47	0/943	0.78	0/1269
45	DO	0.49	0/943	0.80	0/1269
46	BP	0.42	0/1131	0.94	5/1504 (0.3%)
46	DP	0.40	0/1131	0.94	5/1504 (0.3%)
47	BQ	0.41	0/1100	0.76	0/1470
47	DQ	0.42	0/1100	0.78	0/1470
48	BR	0.39	0/974	0.75	0/1302
48	DR	0.42	0/974	0.77	0/1302
49	BS	0.42	0/779	0.71	0/1038
49	DS	0.39	0/779	0.70	0/1038
50	BT	0.44	0/1156	0.77	1/1544 (0.1%)
50	DT	0.45	0/1156	0.77	1/1544 (0.1%)
51	BU	0.40	0/975	0.75	1/1297 (0.1%)
51	DU	0.46	0/975	0.77	1/1297 (0.1%)
52	BV	0.41	0/789	0.74	1/1054 (0.1%)
52	DV	0.42	0/789	0.75	1/1054 (0.1%)
53	BW	0.42	0/907	0.67	0/1216
53	DW	0.42	0/907	0.66	0/1216
54	BX	0.47	0/740	0.88	2/995 (0.2%)
54	DX	0.47	0/740	0.89	2/995 (0.2%)
55	BY	0.43	0/789	0.78	0/1053
55	DY	0.44	0/789	0.77	0/1053
56	BZ	0.40	0/1436	0.74	0/1951
56	DZ	0.40	0/1436	0.77	1/1951 (0.1%)
All	All	0.51	7/313639 (0.0%)	0.74	117/468340 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	AA	0	17
1	CA	0	20
22	CV	0	2
23	AW	0	1
24	AX	0	3
35	BA	1	69
35	DA	1	63
All	All	2	175

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
35	BA	1543	C	N1-C2	5.59	1.45	1.40
40	DF	65	TRP	CB-CG	-5.48	1.40	1.50
35	DA	1543	C	N1-C2	5.43	1.45	1.40
35	DA	652	C	C3'-O3'	5.38	1.49	1.42
35	BA	652	C	C3'-O3'	5.30	1.49	1.42
40	BF	65	TRP	CB-CG	-5.17	1.41	1.50
35	DA	2077	A	C5-C6	-5.07	1.36	1.41

All (117) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
35	DA	2128	C	C2'-C3'-O3'	10.31	132.19	109.50
35	BA	2128	C	C2'-C3'-O3'	10.29	132.14	109.50
46	BP	52	GLU	N-CA-C	8.91	135.05	111.00
46	DP	52	GLU	N-CA-C	8.80	134.76	111.00
38	BD	238	GLY	N-CA-C	-8.62	91.56	113.10
38	DD	238	GLY	N-CA-C	-8.22	92.55	113.10
1	AA	1498	U	C2'-C3'-O3'	7.75	126.55	109.50
46	BP	53	GLY	N-CA-C	-7.69	93.88	113.10
46	DP	53	GLY	N-CA-C	-7.68	93.90	113.10
35	DA	975	C	N1-C1'-C2'	7.41	123.63	114.00
35	BA	283	A	C2'-C3'-O3'	7.28	125.52	109.50
46	BP	59	LEU	CA-CB-CG	7.24	131.95	115.30
35	DA	283	A	C2'-C3'-O3'	7.23	125.40	109.50
35	BA	387	U	C2'-C3'-O3'	7.20	125.33	109.50
35	DA	387	U	C2'-C3'-O3'	7.04	124.99	109.50
28	B2	53	LEU	N-CA-C	-6.87	92.45	111.00
35	BA	975	C	N1-C1'-C2'	6.83	122.88	114.00
27	B1	55	GLY	N-CA-C	-6.63	96.52	113.10
46	DP	59	LEU	CA-CB-CG	6.60	130.48	115.30
1	AA	575	G	N9-C1'-C2'	6.60	122.58	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
54	DX	79	ALA	N-CA-C	-6.41	93.71	111.00
54	BX	79	ALA	N-CA-C	-6.31	93.96	111.00
35	DA	975	C	O4'-C1'-N1	6.24	113.19	108.20
35	BA	975	C	O4'-C1'-N1	6.24	113.19	108.20
35	BA	1616	A	N9-C1'-C2'	6.19	122.05	114.00
35	DA	752	A	C2'-C3'-O3'	6.16	123.55	113.70
35	DA	1616	A	N9-C1'-C2'	6.15	122.00	114.00
27	D1	55	GLY	N-CA-C	-6.13	97.76	113.10
35	BA	2702	U	N1-C1'-C2'	6.06	121.88	114.00
37	BC	140	PRO	N-CA-CB	6.06	110.57	103.30
56	DZ	110	GLY	N-CA-C	-6.02	98.04	113.10
27	D1	37	ILE	CB-CA-C	-6.01	99.58	111.60
37	DC	140	PRO	N-CA-CB	5.99	110.49	103.30
1	CA	575	G	N9-C1'-C2'	5.98	121.78	114.00
39	BE	186	GLY	N-CA-C	5.91	127.88	113.10
35	BA	752	A	C2'-C3'-O3'	5.89	123.13	113.70
38	BD	237	GLU	N-CA-C	5.87	126.86	111.00
34	B8	33	ASN	N-CA-C	-5.84	95.24	111.00
39	DE	186	GLY	N-CA-C	5.83	127.68	113.10
35	BA	669	G	C2'-C3'-O3'	5.83	123.03	113.70
37	BC	181	PRO	N-CA-CB	5.83	110.29	103.30
30	B4	29	PRO	N-CA-CB	5.79	110.25	103.30
35	DA	2702	U	N1-C1'-C2'	5.75	121.47	114.00
1	AA	1225	A	N9-C1'-C2'	5.73	121.45	114.00
37	BC	201	PRO	N-CA-CB	5.73	110.17	103.30
35	DA	178	G	N9-C1'-C2'	-5.73	105.70	112.00
35	BA	193	U	C5'-C4'-C3'	-5.71	106.87	116.00
1	CA	1054	C	N1-C1'-C2'	5.71	121.42	114.00
30	D4	11	PRO	N-CA-CB	5.67	110.11	103.30
30	D4	29	PRO	N-CA-CB	5.67	110.11	103.30
28	D2	55	ARG	N-CA-C	-5.66	95.73	111.00
35	BA	1340	U	N1-C1'-C2'	5.65	121.34	114.00
41	BG	54	GLU	N-CA-C	-5.63	95.80	111.00
37	DC	201	PRO	N-CA-CB	5.63	110.06	103.30
37	DC	181	PRO	N-CA-CB	5.62	110.04	103.30
37	BC	182	PRO	N-CA-CB	5.61	110.04	103.30
34	D8	33	ASN	N-CA-C	-5.57	95.96	111.00
35	BA	254	G	N9-C1'-C2'	-5.55	105.89	112.00
27	D1	64	ALA	N-CA-C	-5.55	96.02	111.00
37	DC	174	PRO	N-CA-CB	5.55	109.96	103.30
37	DC	182	PRO	N-CA-CB	5.54	109.94	103.30
38	DD	237	GLU	N-CA-C	5.52	125.91	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
37	DC	133	PRO	N-CA-CB	5.49	109.89	103.30
1	CA	1529	G	N9-C1'-C2'	5.49	121.13	114.00
37	BC	174	PRO	N-CA-CB	5.48	109.87	103.30
46	BP	48	PRO	N-CA-C	-5.47	97.88	112.10
46	DP	48	PRO	N-CA-C	-5.44	97.95	112.10
37	BC	133	PRO	N-CA-CB	5.42	109.81	103.30
42	DH	156	ALA	N-CA-C	-5.42	96.36	111.00
30	B4	11	PRO	N-CA-CB	5.42	109.80	103.30
35	DA	669	G	C2'-C3'-O3'	5.41	122.36	113.70
41	BG	125	PHE	N-CA-C	-5.41	96.40	111.00
52	DV	18	LEU	CA-CB-CG	5.39	127.69	115.30
33	D7	4	THR	N-CA-C	5.38	125.53	111.00
1	AA	1054	C	N1-C1'-C2'	5.38	120.99	114.00
1	AA	1067	A	C2'-C3'-O3'	5.36	122.27	113.70
30	B4	41	PRO	N-CA-CB	5.35	109.72	103.30
42	BH	156	ALA	N-CA-C	-5.35	96.55	111.00
50	DT	11	GLU	N-CA-C	-5.34	96.58	111.00
37	DC	220	PRO	N-CA-CB	5.33	109.70	103.30
35	BA	178	G	N9-C1'-C2'	-5.33	106.14	112.00
37	BC	220	PRO	N-CA-CB	5.31	109.67	103.30
35	DA	193	U	C5'-C4'-C3'	-5.30	107.52	116.00
54	DX	64	LYS	N-CA-C	5.28	125.25	111.00
1	AA	575	G	C2'-C3'-O3'	5.27	122.13	113.70
1	CA	1225	A	N9-C1'-C2'	5.24	120.81	114.00
54	BX	64	LYS	N-CA-C	5.22	125.11	111.00
22	CV	42	C	C2'-C3'-O3'	5.21	122.03	113.70
35	DA	1493	C	N1-C1'-C2'	5.21	120.77	114.00
51	DU	7	GLY	N-CA-C	5.20	126.11	113.10
52	BV	18	LEU	CA-CB-CG	5.20	127.26	115.30
35	BA	1493	C	N1-C1'-C2'	5.18	120.74	114.00
46	DP	54	GLY	N-CA-C	-5.18	100.15	113.10
1	AA	250	A	C2'-C3'-O3'	5.16	121.96	113.70
35	BA	2422	A	C2'-C3'-O3'	5.15	121.94	113.70
30	D4	7	PRO	N-CA-CB	5.15	109.48	103.30
35	BA	258	G	N9-C1'-C2'	-5.14	106.34	112.00
50	BT	11	GLU	N-CA-C	-5.14	97.12	111.00
30	D4	41	PRO	N-CA-CB	5.14	109.46	103.30
1	CA	1064	G	C2'-C3'-O3'	5.13	121.91	113.70
46	BP	54	GLY	N-CA-C	-5.13	100.27	113.10
35	BA	1396	U	N1-C1'-C2'	5.12	120.66	114.00
35	DA	254	G	N9-C1'-C2'	-5.12	106.37	112.00
35	BA	2524	G	C5'-C4'-C3'	-5.11	107.83	116.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
35	DA	1674	G	N9-C1'-C2'	-5.10	106.39	112.00
35	DA	2524	G	C5'-C4'-C3'	-5.09	107.86	116.00
1	AA	1502	A	N9-C1'-C2'	5.09	120.61	114.00
1	AA	1064	G	C2'-C3'-O3'	5.07	121.81	113.70
35	BA	49	A	N9-C1'-C2'	5.06	120.58	114.00
35	DA	1053	C	N1-C1'-C2'	5.06	120.58	114.00
35	DA	49	A	N9-C1'-C2'	5.06	120.58	114.00
28	B2	55	ARG	N-CA-C	-5.05	97.36	111.00
35	DA	614(C)	A	C2'-C3'-O3'	5.05	121.78	113.70
1	AA	484	G	N9-C1'-C2'	5.03	120.54	114.00
1	AA	115	G	N9-C1'-C2'	5.03	120.54	114.00
35	DA	2897	U	C2'-C3'-O3'	5.01	121.72	113.70
51	BU	7	GLY	N-CA-C	5.01	125.63	113.10

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
35	BA	2128	C	C3'
35	DA	2128	C	C3'

All (175) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AA	1397	C	Sidechain
1	AA	1431	C	Sidechain
1	AA	1433	A	Sidechain
1	AA	21	G	Sidechain
1	AA	265	G	Sidechain
1	AA	323	U	Sidechain
1	AA	445	G	Sidechain
1	AA	587	G	Sidechain
1	AA	626	U	Sidechain
1	AA	639	G	Sidechain
1	AA	686	U	Sidechain
1	AA	741	G	Sidechain
1	AA	792	A	Sidechain
1	AA	832	C	Sidechain
1	AA	884	U	Sidechain
1	AA	916	G	Sidechain
1	AA	96	U	Sidechain
23	AW	74	A	Sidechain
24	AX	16	U	Sidechain

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Mol	Chain	Res	Type	Group
24	AX	17	U	Sidechain
24	AX	22	U	Sidechain
35	BA	102	G	Sidechain
35	BA	1190	G	Sidechain
35	BA	1192	G	Sidechain
35	BA	122	G	Sidechain
35	BA	1294	U	Sidechain
35	BA	1302	A	Sidechain
35	BA	1326	U	Sidechain
35	BA	135	G	Sidechain
35	BA	1357	U	Sidechain
35	BA	1379	A	Sidechain
35	BA	1560	G	Sidechain
35	BA	1613	G	Sidechain
35	BA	1615	C	Sidechain
35	BA	1617	C	Sidechain
35	BA	1647	G	Sidechain
35	BA	1667	G	Sidechain
35	BA	1673	U	Sidechain
35	BA	1674	G	Sidechain
35	BA	1692	U	Sidechain
35	BA	1767	C	Sidechain
35	BA	1772	G	Sidechain
35	BA	1775	U	Sidechain
35	BA	1783	A	Sidechain
35	BA	1806	C	Sidechain
35	BA	1820	U	Sidechain
35	BA	1822	G	Sidechain
35	BA	1834	U	Sidechain
35	BA	1955	U	Sidechain
35	BA	197	A	Sidechain
35	BA	1970	A	Sidechain
35	BA	1980	G	Sidechain
35	BA	1995	U	Sidechain
35	BA	2009	G	Sidechain
35	BA	2031	A	Sidechain
35	BA	2059	A	Sidechain
35	BA	2061	G	Sidechain
35	BA	2069	G	Sidechain
35	BA	2079	U	Sidechain
35	BA	2086	U	Sidechain
35	BA	212	G	Sidechain

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Mol	Chain	Res	Type	Group
35	BA	2228	G	Sidechain
35	BA	2320	A	Sidechain
35	BA	2427	C	Sidechain
35	BA	243	U	Sidechain
35	BA	2437	U	Sidechain
35	BA	2443	C	Sidechain
35	BA	2448	A	Sidechain
35	BA	2517	C	Sidechain
35	BA	2518	A	Sidechain
35	BA	2582	G	Sidechain
35	BA	2608	G	Sidechain
35	BA	2702	U	Sidechain
35	BA	283	A	Sidechain
35	BA	49	A	Sidechain
35	BA	497	A	Sidechain
35	BA	532	A	Sidechain
35	BA	562	U	Sidechain
35	BA	588	U	Sidechain
35	BA	70	G	Sidechain
35	BA	726	G	Sidechain
35	BA	729	G	Sidechain
35	BA	767	U	Sidechain
35	BA	781	A	Sidechain
35	BA	860	U	Sidechain
35	BA	911	A	Sidechain
35	BA	938	G	Sidechain
35	BA	945	A	Sidechain
35	BA	959	A	Sidechain
35	BA	963	U	Sidechain
1	CA	1018	C	Sidechain
1	CA	1414	U	Sidechain
1	CA	1472	U	Sidechain
1	CA	1485	U	Sidechain
1	CA	1498	U	Sidechain
1	CA	21	G	Sidechain
1	CA	265	G	Sidechain
1	CA	300	A	Sidechain
1	CA	323	U	Sidechain
1	CA	575	G	Sidechain
1	CA	587	G	Sidechain
1	CA	639	G	Sidechain
1	CA	686	U	Sidechain

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Mol	Chain	Res	Type	Group
1	CA	741	G	Sidechain
1	CA	783	C	Sidechain
1	CA	814	A	Sidechain
1	CA	832	C	Sidechain
1	CA	884	U	Sidechain
1	CA	916	G	Sidechain
1	CA	96	U	Sidechain
22	CV	29	G	Sidechain
22	CV	32	U	Sidechain
35	DA	1158	C	Sidechain
35	DA	1192	G	Sidechain
35	DA	1268	A	Sidechain
35	DA	1302	A	Sidechain
35	DA	1326	U	Sidechain
35	DA	135	G	Sidechain
35	DA	1357	U	Sidechain
35	DA	1379	A	Sidechain
35	DA	1411	C	Sidechain
35	DA	1560	G	Sidechain
35	DA	1608	A	Sidechain
35	DA	1613	G	Sidechain
35	DA	1615	C	Sidechain
35	DA	1617	C	Sidechain
35	DA	1647	G	Sidechain
35	DA	1673	U	Sidechain
35	DA	1674	G	Sidechain
35	DA	1767	C	Sidechain
35	DA	1772	G	Sidechain
35	DA	1777	U	Sidechain
35	DA	1783	A	Sidechain
35	DA	1786	A	Sidechain
35	DA	1806	C	Sidechain
35	DA	1820	U	Sidechain
35	DA	1822	G	Sidechain
35	DA	1834	U	Sidechain
35	DA	1955	U	Sidechain
35	DA	197	A	Sidechain
35	DA	1980	G	Sidechain
35	DA	1985	G	Sidechain
35	DA	2009	G	Sidechain
35	DA	2010	G	Sidechain
35	DA	2031	A	Sidechain

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Mol	Chain	Res	Type	Group
35	DA	204	A	Sidechain
35	DA	2052	G	Sidechain
35	DA	2059	A	Sidechain
35	DA	2061	G	Sidechain
35	DA	2079	U	Sidechain
35	DA	2083	G	Sidechain
35	DA	2086	U	Sidechain
35	DA	2396	G	Sidechain
35	DA	2427	C	Sidechain
35	DA	243	U	Sidechain
35	DA	2437	U	Sidechain
35	DA	2443	C	Sidechain
35	DA	2518	A	Sidechain
35	DA	2564	A	Sidechain
35	DA	2689	U	Sidechain
35	DA	2702	U	Sidechain
35	DA	49	A	Sidechain
35	DA	497	A	Sidechain
35	DA	532	A	Sidechain
35	DA	562	U	Sidechain
35	DA	588	U	Sidechain
35	DA	70	G	Sidechain
35	DA	726	G	Sidechain
35	DA	767	U	Sidechain
35	DA	787	U	Sidechain
35	DA	835	A	Sidechain
35	DA	860	U	Sidechain
35	DA	911	A	Sidechain
35	DA	945	A	Sidechain
35	DA	963	U	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AA	32329	0	16313	2049	0
1	CA	32329	0	16314	2040	0
2	AB	1901	0	1951	382	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	CB	1901	0	1951	378	0
3	AC	1613	0	1677	256	0
3	CC	1613	0	1677	246	0
4	AD	1703	0	1763	346	0
4	CD	1703	0	1764	336	0
5	AE	1147	0	1207	211	0
5	CE	1147	0	1207	202	0
6	AF	843	0	857	151	0
6	CF	843	0	857	157	0
7	AG	1257	0	1296	235	0
7	CG	1257	0	1296	216	0
8	AH	1116	0	1177	231	0
8	CH	1116	0	1177	233	0
9	AI	1011	0	1041	208	0
9	CI	1011	0	1041	209	0
10	AJ	795	0	840	153	0
10	CJ	795	0	840	154	0
11	AK	885	0	904	179	0
11	CK	885	0	904	180	0
12	AL	971	0	1057	210	0
12	CL	971	0	1057	196	0
13	AM	988	0	1055	163	0
13	CM	988	0	1055	156	0
14	AN	492	0	529	95	0
14	CN	492	0	530	89	0
15	AO	734	0	771	111	0
15	CO	734	0	771	117	0
16	AP	701	0	720	157	0
16	CP	701	0	720	146	0
17	AQ	824	0	891	140	0
17	CQ	824	0	891	138	0
18	AR	574	0	644	138	0
18	CR	574	0	644	133	0
19	AS	630	0	652	139	0
19	CS	630	0	652	138	0
20	AT	763	0	861	133	0
20	CT	763	0	861	125	0
21	AU	209	0	221	30	0
21	CU	209	0	221	28	0
22	AV	362	0	186	23	0
22	CV	362	0	186	20	0
23	AW	1641	0	836	105	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
23	CW	1641	0	839	72	0
24	AX	230	0	118	12	0
24	CX	230	0	117	16	0
25	AY	1478	0	1526	283	0
25	CY	1478	0	1526	303	0
26	B0	652	0	658	102	0
26	D0	652	0	658	87	0
27	B1	693	0	764	248	0
27	D1	693	0	763	241	0
28	B2	421	0	461	118	0
28	D2	421	0	461	146	0
29	B3	468	0	523	86	0
29	D3	468	0	523	82	0
30	B4	242	0	103	14	0
30	D4	242	0	103	10	0
31	B5	459	0	480	103	0
31	D5	459	0	480	107	0
32	B6	381	0	390	69	0
32	D6	381	0	390	72	0
33	B7	419	0	467	75	0
33	D7	419	0	467	72	0
34	B8	508	0	576	154	0
34	D8	508	0	576	143	0
35	BA	59601	0	30029	3750	1
35	DA	59601	0	30026	3756	0
36	BB	2551	0	1294	137	1
36	DB	2551	0	1294	155	0
37	BC	1142	0	861	85	0
37	DC	1143	0	865	78	0
38	BD	2105	0	2182	497	0
38	DD	2105	0	2182	503	0
39	BE	1564	0	1629	382	0
39	DE	1564	0	1629	393	0
40	BF	1624	0	1677	349	0
40	DF	1624	0	1676	353	0
41	BG	1474	0	1534	336	0
41	DG	1474	0	1534	441	0
42	BH	1223	0	1282	242	0
42	DH	1223	0	1282	245	0
43	BI	1132	0	1218	227	0
43	DI	1132	0	1218	225	0
44	BN	1105	0	1180	277	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
44	DN	1105	0	1180	291	0
45	BO	933	0	996	234	0
45	DO	933	0	996	237	0
46	BP	1114	0	1187	382	0
46	DP	1114	0	1187	378	0
47	BQ	1080	0	1127	274	0
47	DQ	1080	0	1127	259	0
48	BR	960	0	1021	221	0
48	DR	960	0	1021	220	0
49	BS	771	0	832	201	0
49	DS	771	0	832	207	0
50	BT	1142	0	1202	320	0
50	DT	1142	0	1202	325	0
51	BU	958	0	1014	245	0
51	DU	958	0	1014	256	0
52	BV	779	0	851	231	0
52	DV	779	0	851	234	0
53	BW	896	0	953	174	0
53	DW	896	0	953	164	0
54	BX	726	0	778	251	0
54	DX	726	0	777	252	0
55	BY	776	0	870	205	0
55	DY	776	0	870	208	0
56	BZ	1404	0	1432	353	0
56	DZ	1404	0	1432	318	0
57	AA	204	0	0	0	0
57	AD	2	0	0	0	0
57	AE	2	0	0	0	0
57	AG	1	0	0	0	0
57	AK	1	0	0	0	0
57	AL	2	0	0	0	0
57	AM	1	0	0	0	0
57	AV	4	0	0	0	0
57	AW	23	0	0	0	0
57	AX	5	0	0	0	0
57	B1	4	0	0	0	0
57	B2	2	0	0	0	0
57	B3	1	0	0	0	0
57	B5	2	0	0	0	0
57	B7	1	0	0	0	0
57	BA	445	0	0	1	0
57	BB	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
57	BD	3	0	0	0	0
57	BE	1	0	0	0	0
57	BF	4	0	0	0	0
57	BG	1	0	0	0	0
57	BN	4	0	0	0	0
57	BO	1	0	0	0	0
57	BP	3	0	0	0	0
57	BU	2	0	0	1	0
57	BX	2	0	0	0	0
57	CA	208	0	0	0	0
57	CE	3	0	0	0	0
57	CG	1	0	0	0	0
57	CL	2	0	0	0	0
57	CN	1	0	0	0	0
57	CV	4	0	0	0	0
57	CW	23	0	0	0	0
57	CX	4	0	0	0	0
57	D1	4	0	0	0	0
57	D2	2	0	0	0	0
57	D3	1	0	0	0	0
57	D5	1	0	0	0	0
57	D7	1	0	0	0	0
57	DA	441	0	0	0	0
57	DB	19	0	0	0	0
57	DD	3	0	0	0	0
57	DE	2	0	0	0	0
57	DF	5	0	0	0	0
57	DG	1	0	0	0	0
57	DH	1	0	0	0	0
57	DN	2	0	0	0	0
57	DP	4	0	0	0	0
57	DS	1	0	0	0	0
57	DU	3	0	0	0	0
57	DV	1	0	0	0	0
57	DX	3	0	0	0	0
58	AD	1	0	0	0	0
58	AN	1	0	0	0	0
58	CD	1	0	0	0	0
58	CN	1	0	0	0	0
All	All	290487	0	197331	30322	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 62.

All (30322) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:29:LEU:HD21	53:BW:33:ARG:HE	1.07	1.20
55:DY:45:VAL:HA	55:DY:62:GLU:HG2	1.20	1.20
35:DA:2758:A:H2'	35:DA:2759:G:H5''	1.25	1.19
1:CA:1442(A):G:H3'	1:CA:1442(B):A:H5''	1.22	1.18
43:BI:91:SER:HB2	43:BI:119:PRO:HB2	1.21	1.18
1:CA:979:C:H3'	1:CA:980:C:H5''	1.22	1.18
53:DW:29:LEU:HD21	53:DW:33:ARG:HE	1.08	1.17
38:BD:27:THR:HG23	38:BD:28:GLU:H	1.09	1.17
41:DG:5:VAL:HG12	41:DG:6:ALA:H	1.07	1.17
1:CA:1397:C:H42	24:CX:22:U:H3'	1.02	1.16
35:DA:1798:U:H5'	38:DD:259:THR:HG22	1.16	1.16
16:CP:20:VAL:HG21	16:CP:32:TYR:HB2	1.27	1.16
35:BA:2758:A:H2'	35:BA:2759:G:H5''	1.26	1.15
38:BD:8:PRO:HB3	38:BD:14:ARG:HD3	1.26	1.15
41:DG:111:LEU:HB2	41:DG:112:PRO:HD3	1.25	1.15
1:CA:1123:A:H4'	10:CJ:36:GLY:HA3	1.24	1.14
34:B8:46:ARG:HB3	34:B8:46:ARG:HH11	1.03	1.14
47:DQ:39:PRO:HB3	47:DQ:99:PRO:HD3	1.20	1.14
50:BT:80:SER:HB3	50:BT:81:PRO:CD	1.78	1.13
16:AP:20:VAL:HG21	16:AP:32:TYR:HB2	1.26	1.13
36:BB:74:U:H2'	36:BB:75:G:H5''	1.27	1.13
35:DA:2491:U:H5'	35:DA:2570:G:H5''	1.30	1.13
35:BA:2893:G:H5'	35:BA:2894:G:H5'	1.23	1.13
1:AA:1189:C:H5''	3:AC:5:ILE:HG21	1.30	1.13
35:DA:1494:A:H2'	35:DA:1495:A:H5''	1.16	1.13
52:BV:28:GLU:HB2	52:BV:29:PRO:HD3	1.24	1.12
42:DH:102:ALA:HB2	42:DH:117:PRO:HD3	1.31	1.12
1:AA:979:C:H3'	1:AA:980:C:H5''	1.23	1.12
34:D8:62:LEU:HD13	35:DA:242:G:H5''	1.25	1.12
54:BX:72:LYS:HG3	54:BX:74:PRO:HD3	1.16	1.12
10:AJ:50:ILE:H	10:AJ:50:ILE:HD13	1.15	1.12
35:DA:2893:G:H5'	35:DA:2894:G:H5'	1.20	1.12
50:DT:80:SER:HB3	50:DT:81:PRO:CD	1.77	1.12
56:BZ:3:TYR:N	56:BZ:57:ILE:HA	1.64	1.12
35:DA:1879:C:H2'	35:DA:1880:C:H5''	1.30	1.11
27:B1:58:ILE:HD13	27:B1:59:THR:H	1.06	1.11
2:AB:75:LYS:HA	2:AB:78:GLN:HE21	1.12	1.11
38:DD:161:THR:OG1	38:DD:196:VAL:HG21	1.51	1.11
2:AB:96:ARG:H	2:AB:96:ARG:HD2	1.02	1.11
47:BQ:140:ALA:HB3	56:BZ:53:ILE:HD12	1.27	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:60:LEU:HD22	41:BG:63:ILE:HD11	1.25	1.11
41:DG:114:ILE:HG21	41:DG:117:PHE:HB2	1.27	1.11
50:BT:62:THR:HG22	50:BT:75:ILE:HA	1.32	1.11
55:DY:95:LYS:HG2	55:DY:100:ALA:HA	1.29	1.11
34:B8:62:LEU:HD13	35:BA:242:G:H5''	1.28	1.11
28:D2:44:LEU:HD23	35:DA:61:G:H5'	1.32	1.11
32:D6:10:LEU:HD12	34:D8:36:LYS:HD3	1.27	1.10
10:CJ:50:ILE:HD13	10:CJ:50:ILE:H	1.14	1.10
35:BA:2598:A:H5''	38:BD:236:GLY:H	1.05	1.10
41:BG:27:ASN:HD21	41:BG:29:TRP:HB2	1.16	1.10
42:BH:102:ALA:HB2	42:BH:117:PRO:HD3	1.30	1.10
35:BA:483:A:H1'	55:BY:47:LYS:HG2	1.32	1.10
41:DG:36:LYS:HD3	41:DG:95:ARG:HH22	1.14	1.10
35:BA:2639:A:H2'	35:BA:2640:G:H5''	1.34	1.10
39:BE:51:PHE:H	39:BE:74:PRO:HG3	1.17	1.10
56:DZ:166:SER:CB	56:DZ:168:GLU:H	1.64	1.10
52:DV:28:GLU:HB2	52:DV:29:PRO:HD3	1.25	1.10
35:BA:1494:A:H2'	35:BA:1495:A:H5''	1.18	1.09
35:DA:1884:A:H2'	35:DA:1885:A:H5''	1.33	1.09
35:DA:2598:A:H5''	38:DD:236:GLY:H	1.07	1.09
1:AA:1123:A:H4'	10:AJ:36:GLY:HA3	1.22	1.09
14:AN:4:LYS:HD2	14:AN:7:ILE:HD11	1.32	1.09
32:B6:10:LEU:HD12	34:B8:36:LYS:HD3	1.21	1.09
1:CA:1189:C:H5''	3:CC:5:ILE:HG21	1.33	1.09
38:BD:35:LYS:HD3	38:BD:63:ARG:HB3	1.31	1.09
38:DD:8:PRO:HB3	38:DD:14:ARG:HD3	1.25	1.09
41:DG:51:ARG:NE	41:DG:51:ARG:HA	1.63	1.09
56:DZ:166:SER:HB2	56:DZ:168:GLU:H	1.14	1.09
3:CC:43:LEU:HD22	3:CC:47:LEU:HD22	1.34	1.09
38:DD:27:THR:HG23	38:DD:28:GLU:H	1.10	1.09
47:DQ:52:VAL:HG13	47:DQ:53:ALA:H	1.17	1.09
35:BA:1879:C:H2'	35:BA:1880:C:H5''	1.28	1.08
1:CA:1321:C:H5'	1:CA:1322:C:H5''	1.35	1.08
45:DO:114:ILE:HD12	45:DO:114:ILE:H	1.18	1.08
32:D6:51:GLU:HG2	32:D6:52:VAL:H	1.18	1.08
32:B6:51:GLU:HG2	32:B6:52:VAL:H	1.18	1.08
49:BS:15:ARG:HB3	49:BS:18:ILE:HD13	1.32	1.08
4:CD:22:LYS:HB2	4:CD:26:CYS:HB2	1.34	1.08
36:DB:74:U:H2'	36:DB:75:G:H5''	1.29	1.08
43:DI:91:SER:HB2	43:DI:119:PRO:HB2	1.18	1.08
50:DT:109:GLU:HB3	50:DT:113:LYS:HE3	1.36	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BQ:39:PRO:HB3	47:BQ:99:PRO:HD3	1.18	1.08
35:DA:2639:A:H2'	35:DA:2640:G:H5''	1.32	1.08
27:B1:42:GLN:HG2	27:B1:43:TYR:H	1.15	1.08
35:BA:2284:C:H2'	35:BA:2285:C:H5''	1.29	1.08
44:BN:66:LYS:HD3	44:BN:70:LYS:HB2	1.29	1.08
35:DA:2284:C:H2'	35:DA:2285:C:H5''	1.32	1.08
35:DA:483:A:H1'	55:DY:47:LYS:HG2	1.34	1.08
3:AC:9:GLY:HA2	3:AC:12:LEU:HD23	1.36	1.07
25:AY:39:LEU:HB2	25:AY:53:ASN:HB3	1.30	1.07
35:BA:2491:U:H5'	35:BA:2570:G:H5''	1.32	1.07
40:BF:2:LYS:HG3	40:BF:25:PRO:HB2	1.30	1.07
55:BY:45:VAL:HA	55:BY:62:GLU:HG2	1.23	1.07
5:CE:50:GLU:HG3	5:CE:52:PRO:HD2	1.34	1.07
14:CN:4:LYS:HD2	14:CN:7:ILE:HD11	1.30	1.07
46:DP:71:VAL:CG1	46:DP:72:PRO:HD3	1.83	1.07
27:D1:10:LYS:HG3	27:D1:11:ARG:H	1.18	1.07
34:D8:46:ARG:HH11	34:D8:46:ARG:HB3	1.02	1.07
35:BA:271(P):C:H5'	43:BI:46:ALA:HB2	1.32	1.07
44:DN:66:LYS:HD3	44:DN:70:LYS:HB2	1.28	1.07
34:B8:52:LYS:H	34:B8:53:PRO:HD2	1.19	1.07
35:BA:1798:U:H5'	38:BD:259:THR:HG22	1.14	1.07
46:BP:71:VAL:CG1	46:BP:72:PRO:HD3	1.83	1.07
3:CC:9:GLY:HA2	3:CC:12:LEU:HD23	1.36	1.07
43:BI:77:LEU:HB2	43:BI:140:LEU:HD13	1.36	1.07
40:DF:2:LYS:HG3	40:DF:25:PRO:HB2	1.30	1.07
35:BA:1586:A:H3'	35:BA:1587:A:H5''	1.07	1.07
55:BY:95:LYS:HG2	55:BY:100:ALA:HA	1.27	1.07
2:CB:96:ARG:H	2:CB:96:ARG:HD2	1.03	1.07
47:BQ:75:THR:HA	47:BQ:88:GLY:HA3	1.37	1.07
35:DA:1586:A:H3'	35:DA:1587:A:H5''	1.10	1.07
35:DA:271(P):C:H5'	43:DI:46:ALA:HB2	1.33	1.07
35:BA:2821:A:P	48:BR:2:ARG:HH22	1.77	1.06
41:DG:51:ARG:NH1	41:DG:53:LEU:HG	1.70	1.06
35:DA:2821:A:P	48:DR:2:ARG:HH22	1.77	1.06
3:AC:43:LEU:HD22	3:AC:47:LEU:HD22	1.33	1.06
38:DD:25:THR:HG21	38:DD:81:ALA:HB1	1.35	1.06
38:DD:35:LYS:HD3	38:DD:63:ARG:HB3	1.34	1.06
5:AE:50:GLU:HG3	5:AE:52:PRO:HD2	1.37	1.06
35:BA:1286:A:H2'	35:BA:1288:U:OP2	1.54	1.06
52:BV:83:ARG:HG2	52:BV:83:ARG:HH11	1.16	1.06
28:D2:41:ILE:HD12	28:D2:41:ILE:H	1.21	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:62:THR:HG22	50:DT:75:ILE:HA	1.31	1.06
50:BT:66:VAL:HA	50:BT:71:GLY:HA2	1.38	1.05
48:DR:10:LEU:HD22	48:DR:17:ARG:HD2	1.36	1.05
35:BA:2562:U:H1'	45:BO:23:ARG:HH12	1.21	1.05
1:CA:908:A:H2'	1:CA:909:A:H8	1.18	1.05
38:BD:161:THR:OG1	38:BD:196:VAL:HG21	1.56	1.05
49:DS:15:ARG:HB3	49:DS:18:ILE:HD13	1.29	1.05
26:B0:70:GLN:HG2	26:B0:71:ASP:H	1.20	1.05
28:B2:30:ARG:H	28:B2:30:ARG:HD2	1.18	1.05
35:DA:782:A:C2	38:DD:226:MET:HG2	1.92	1.05
45:BO:63:VAL:HG22	45:BO:84:ALA:HA	1.38	1.05
2:CB:75:LYS:HA	2:CB:78:GLN:HE21	1.12	1.05
43:BI:72:LEU:HD12	43:BI:138:ILE:HD11	1.38	1.05
4:CD:120:LEU:HD12	4:CD:120:LEU:H	1.18	1.05
51:DU:92:ARG:HB3	52:DV:11:GLN:NE2	1.70	1.05
2:AB:71:VAL:HG22	2:AB:93:VAL:H	1.21	1.05
35:DA:2533:A:H2'	35:DA:2534:A:H5''	1.36	1.05
50:DT:28:VAL:HG21	50:DT:47:GLY:H	1.20	1.05
39:BE:179:GLU:HB3	39:BE:181:LEU:HD21	1.37	1.04
27:D1:33:LYS:HG2	27:D1:34:THR:H	1.16	1.04
39:DE:108:SER:HB3	39:DE:165:VAL:HG21	1.38	1.04
44:DN:34:LEU:HD21	44:DN:120:LEU:HB2	1.39	1.04
35:BA:1747(A):G:H2'	35:BA:1748:G:H5''	1.37	1.04
35:BA:2476:A:H2'	35:BA:2477:C:H5''	1.40	1.04
44:BN:34:LEU:HD21	44:BN:120:LEU:HB2	1.37	1.04
45:BO:114:ILE:HD12	45:BO:114:ILE:H	1.20	1.04
27:D1:11:ARG:O	27:D1:13:ILE:N	1.89	1.04
2:CB:71:VAL:HG22	2:CB:93:VAL:H	1.18	1.04
18:CR:37:VAL:HG23	18:CR:38:GLU:H	1.22	1.04
35:BA:2729:G:H1'	39:BE:187:ALA:HB2	1.34	1.04
52:DV:38:LEU:HD23	52:DV:39:LEU:H	1.20	1.04
34:B8:13:ARG:HB3	46:BP:63:PRO:HA	1.37	1.04
35:DA:586:A:H5'	40:DF:89:VAL:HG21	1.37	1.04
27:D1:58:ILE:HD12	27:D1:59:THR:H	1.22	1.04
51:BU:66:ASN:HD21	51:BU:70:ARG:NH2	1.54	1.04
35:DA:689:A:H2'	35:DA:690:G:H8	1.22	1.04
35:DA:2729:G:H1'	39:DE:187:ALA:HB2	1.38	1.04
35:DA:1203:G:H4'	46:DP:7:ARG:HG2	1.36	1.04
56:DZ:70:LEU:HD12	56:DZ:91:LEU:HD21	1.38	1.04
35:BA:1884:A:H2'	35:BA:1885:A:H5''	1.34	1.03
35:BA:782:A:C2	38:BD:226:MET:HG2	1.93	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:86:SER:HA	27:B1:89:GLU:HG3	1.35	1.03
38:BD:27:THR:HG21	38:BD:83:GLU:HG2	1.40	1.03
4:AD:96:LEU:H	4:AD:96:LEU:HD22	1.19	1.03
27:B1:16:ASN:HB3	27:B1:46:LEU:HG	1.39	1.03
38:BD:25:THR:HG21	38:BD:81:ALA:HB1	1.38	1.03
3:CC:15:THR:HG22	3:CC:16:ARG:HH12	1.21	1.03
1:CA:975:A:H4'	1:CA:976:G:H5''	1.38	1.03
35:DA:2476:A:H2'	35:DA:2477:C:H5''	1.40	1.03
46:DP:146:VAL:HG22	46:DP:147:LEU:H	1.24	1.03
48:DR:48:VAL:HA	48:DR:51:LEU:HD12	1.41	1.03
1:AA:1412:C:H2'	1:AA:1413:A:C8	1.94	1.03
40:BF:53:THR:HG23	40:BF:56:GLU:HB2	1.40	1.03
50:DT:66:VAL:HA	50:DT:71:GLY:HA2	1.39	1.03
56:DZ:166:SER:HB2	56:DZ:168:GLU:N	1.74	1.03
3:AC:15:THR:HG22	3:AC:16:ARG:HH12	1.18	1.03
28:D2:49:LYS:HB3	28:D2:53:LEU:HD22	1.34	1.03
35:DA:1286:A:H2'	35:DA:1288:U:OP2	1.58	1.03
38:DD:92:ILE:HG22	38:DD:106:ILE:HA	1.40	1.03
1:AA:1321:C:H5'	1:AA:1322:C:H5''	1.36	1.03
35:BA:2533:A:H2'	35:BA:2534:A:H5''	1.40	1.02
35:BA:2598:A:H5''	38:BD:236:GLY:N	1.73	1.02
35:DA:2348:U:H2'	35:DA:2349:G:H5''	1.41	1.02
34:D8:13:ARG:HB3	46:DP:63:PRO:HA	1.36	1.02
55:DY:10:GLY:HA2	55:DY:27:VAL:HG13	1.40	1.02
1:AA:908:A:H2'	1:AA:909:A:H8	1.20	1.02
1:AA:975:A:H4'	1:AA:976:G:H5''	1.41	1.02
1:AA:585:G:H4'	12:AL:8:ASN:HD21	1.23	1.02
41:DG:41:GLN:HG2	41:DG:43:LEU:HD12	1.37	1.02
4:AD:22:LYS:HB2	4:AD:26:CYS:HB2	1.37	1.02
2:CB:165:VAL:HG23	2:CB:166:ASP:H	1.23	1.02
3:CC:157:ILE:HB	3:CC:164:ARG:HH12	1.23	1.02
28:D2:14:ARG:HG2	28:D2:15:LYS:H	1.20	1.02
35:BA:404:C:H4'	35:BA:405:U:H5'	1.41	1.02
39:DE:51:PHE:H	39:DE:74:PRO:HG3	1.21	1.02
46:DP:38:GLN:HG3	46:DP:39:LYS:H	1.23	1.02
48:DR:98:LEU:HB2	48:DR:113:LEU:HD23	1.41	1.02
40:DF:67:GLN:O	40:DF:67:GLN:HG3	1.56	1.02
2:AB:165:VAL:HG23	2:AB:166:ASP:H	1.24	1.02
48:BR:10:LEU:HD22	48:BR:17:ARG:HD2	1.41	1.02
50:BT:109:GLU:HB3	50:BT:113:LYS:HE3	1.36	1.02
43:DI:72:LEU:HD12	43:DI:138:ILE:HD11	1.39	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:75:THR:HA	47:DQ:88:GLY:HA3	1.37	1.02
44:DN:14:VAL:HG12	44:DN:15:LEU:H	1.25	1.02
56:DZ:56:VAL:HA	56:DZ:70:LEU:HD21	1.04	1.02
32:B6:12:GLU:HA	32:B6:23:THR:HA	1.42	1.02
46:DP:71:VAL:HG13	46:DP:72:PRO:CD	1.90	1.02
39:DE:110:GLY:HA2	39:DE:161:GLY:HA3	1.40	1.02
18:AR:25:THR:HG22	18:AR:42:ARG:HH11	1.22	1.01
35:BA:2348:U:H2'	35:BA:2349:G:H5''	1.38	1.01
35:BA:389:G:H22	46:BP:71:VAL:HG11	1.24	1.01
35:BA:2737:G:H2'	35:BA:2738:A:H8	1.25	1.01
35:BA:389:G:H1	46:BP:71:VAL:HB	1.25	1.01
47:BQ:52:VAL:HG13	47:BQ:53:ALA:H	1.24	1.01
1:CA:1524:C:H2'	1:CA:1525:G:H8	1.25	1.01
18:CR:25:THR:HG22	18:CR:42:ARG:HH11	1.21	1.01
41:BG:5:VAL:HG12	41:BG:6:ALA:H	1.25	1.01
52:BV:18:LEU:HD22	52:BV:19:LYS:H	1.23	1.01
41:DG:132:ASN:HD21	41:DG:157:ILE:HG13	1.23	1.01
53:DW:73:ALA:HB3	53:DW:106:ILE:HD11	1.40	1.01
48:BR:98:LEU:HB2	48:BR:113:LEU:HD23	1.41	1.01
4:AD:120:LEU:H	4:AD:120:LEU:HD12	1.23	1.01
35:BA:94(A):G:H2'	35:BA:95:G:H5''	1.43	1.01
26:D0:70:GLN:HG2	26:D0:71:ASP:H	1.18	1.01
56:DZ:69:THR:HG22	56:DZ:90:VAL:HG22	1.41	1.01
41:BG:173:LEU:HA	41:BG:176:LEU:HD12	1.38	1.01
35:DA:2036:C:H6	35:DA:2036:C:H5'	1.24	1.01
38:DD:226:MET:HE2	38:DD:230:ASP:HB3	1.43	1.01
51:DU:66:ASN:HD21	51:DU:70:ARG:NH2	1.58	1.01
18:AR:37:VAL:HG23	18:AR:38:GLU:H	1.21	1.01
35:DA:2737:G:H2'	35:DA:2738:A:H8	1.25	1.01
42:DH:19:VAL:HB	42:DH:44:VAL:HG13	1.42	1.01
2:AB:112:VAL:O	2:AB:115:LEU:HB3	1.61	1.01
27:B1:27:GLU:HB2	27:B1:33:LYS:O	1.58	1.01
35:DA:94(A):G:H2'	35:DA:95:G:H5''	1.43	1.01
35:DA:342:G:C2'	35:DA:343:C:H5''	1.91	1.01
35:BA:2348:U:C2'	35:BA:2349:G:H5''	1.90	1.01
50:BT:28:VAL:HG21	50:BT:47:GLY:H	1.24	1.01
54:BX:72:LYS:HE3	54:BX:74:PRO:HB3	1.40	1.01
4:CD:96:LEU:H	4:CD:96:LEU:HD22	1.22	1.01
41:DG:132:ASN:HB2	41:DG:159:VAL:HG22	1.41	1.01
16:AP:60:LEU:HD23	16:AP:64:ALA:HB3	1.43	1.00
27:B1:94:LEU:HD22	27:B1:95:LEU:H	1.22	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:146:VAL:HG22	46:BP:147:LEU:H	1.24	1.00
35:DA:1747(A):G:H2'	35:DA:1748:G:H5''	1.38	1.00
38:DD:27:THR:HG21	38:DD:83:GLU:HG2	1.41	1.00
1:AA:939:G:H5''	7:AG:102:ARG:HH22	1.24	1.00
50:BT:11:GLU:H	50:BT:11:GLU:CD	1.60	1.00
51:DU:64:ARG:HH21	51:DU:64:ARG:HB3	1.18	1.00
49:BS:28:VAL:HG12	49:BS:29:PHE:H	1.23	1.00
52:BV:38:LEU:HD23	52:BV:39:LEU:H	1.23	1.00
35:DA:27:G:HO2'	35:DA:28:A:H8	1.02	1.00
46:BP:71:VAL:HG13	46:BP:72:PRO:CD	1.90	1.00
35:BA:1203:G:H4'	46:BP:7:ARG:HG2	1.39	1.00
1:CA:1324:A:H2'	1:CA:1325:C:H6	1.26	1.00
25:CY:10:THR:O	25:CY:14:MET:HG3	1.59	1.00
32:D6:12:GLU:HA	32:D6:23:THR:HA	1.39	1.00
35:DA:1494:A:H2'	35:DA:1495:A:C5'	1.92	1.00
27:B1:10:LYS:O	27:B1:13:ILE:HG22	1.61	1.00
36:DB:74:U:C2'	36:DB:75:G:H5''	1.91	1.00
35:DA:389:G:H22	46:DP:71:VAL:HG11	1.25	1.00
56:DZ:73:GLN:HG2	56:DZ:74:VAL:H	1.25	1.00
39:BE:108:SER:HB3	39:BE:165:VAL:HG21	1.38	1.00
43:BI:38:LEU:HD12	43:BI:38:LEU:H	1.26	1.00
12:CL:55:VAL:HG12	12:CL:56:ALA:H	1.26	1.00
12:CL:83:VAL:HG22	12:CL:84:LEU:H	1.26	1.00
43:DI:77:LEU:HB2	43:DI:140:LEU:HD13	1.40	1.00
35:DA:17:G:HO2'	51:DU:25:TRP:HZ3	1.09	1.00
52:DV:83:ARG:HG2	52:DV:83:ARG:HH11	1.22	1.00
1:AA:1363(A):A:H4'	1:AA:1364:U:H5''	1.40	1.00
35:BA:1586:A:H3'	35:BA:1587:A:C5'	1.91	1.00
1:CA:1216:G:H5''	14:CN:5:ALA:HB2	1.43	1.00
13:CM:91:ARG:HH11	19:CS:81:ARG:HH22	1.06	1.00
1:AA:1324:A:H2'	1:AA:1325:C:H6	1.27	1.00
1:AA:1469:G:H2'	1:AA:1470:G:H8	1.25	1.00
16:AP:26:ARG:HH11	16:AP:26:ARG:HB3	1.26	1.00
35:BA:342:G:C2'	35:BA:343:C:H5''	1.91	1.00
46:BP:71:VAL:HG13	46:BP:72:PRO:HD3	1.00	1.00
35:DA:404:C:H4'	35:DA:405:U:H5'	1.39	1.00
42:DH:18:GLU:HB2	42:DH:25:LYS:HB2	1.43	1.00
38:BD:92:ILE:HG22	38:BD:106:ILE:HA	1.40	1.00
47:BQ:55:VAL:HG12	47:BQ:64:ILE:HD12	1.43	1.00
2:AB:236:TYR:HA	2:AB:239:VAL:HG23	1.44	1.00
19:AS:36:ARG:HH22	19:AS:75:ALA:HB3	1.27	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1348:G:H2'	35:BA:1349:A:H5''	1.42	1.00
35:BA:2498:C:O2'	35:BA:2499:C:H5'	1.62	1.00
44:BN:14:VAL:HG12	44:BN:15:LEU:H	1.25	1.00
1:CA:1452:C:H1'	1:CA:1456:G:N2	1.75	1.00
35:DA:2701:C:H3'	35:DA:2702:U:H5''	1.40	1.00
35:BA:1657:C:H2'	35:BA:1658:C:H6	1.27	0.99
39:BE:110:GLY:HA2	39:BE:161:GLY:HA3	1.44	0.99
16:CP:26:ARG:HH11	16:CP:26:ARG:HB3	1.25	0.99
27:D1:62:VAL:HG21	27:D1:67:ILE:HA	1.39	0.99
52:DV:18:LEU:HD22	52:DV:19:LYS:H	1.26	0.99
34:B8:32:LEU:C	34:B8:34:TRP:H	1.55	0.99
36:BB:74:U:C2'	36:BB:75:G:H5''	1.91	0.99
46:BP:38:GLN:HG3	46:BP:39:LYS:H	1.24	0.99
1:CA:194:C:H2'	1:CA:195:A:H5''	1.44	0.99
25:AY:60:ALA:HA	25:AY:66:LEU:HA	1.41	0.99
52:BV:70:ILE:HB	52:BV:90:PRO:HB2	1.40	0.99
34:D8:52:LYS:H	34:D8:53:PRO:HD2	1.24	0.99
3:AC:157:ILE:HB	3:AC:164:ARG:HH12	1.27	0.99
1:AA:1216:G:H5''	14:AN:5:ALA:HB2	1.45	0.99
27:B1:47:GLN:HG2	35:BA:2230:G:H1'	0.99	0.99
27:B1:85:LEU:C	27:B1:87:PRO:HD3	1.83	0.99
2:CB:236:TYR:HA	2:CB:239:VAL:HG23	1.42	0.99
35:DA:1348:G:H2'	35:DA:1349:A:H5''	1.42	0.99
39:DE:179:GLU:HB3	39:DE:181:LEU:HD21	1.42	0.99
52:DV:64:HIS:HB3	52:DV:96:ILE:HA	1.43	0.99
7:AG:75:VAL:HG12	7:AG:88:PRO:HB3	1.44	0.99
25:AY:133:ARG:HD3	25:AY:165:THR:HG21	1.40	0.99
53:BW:73:ALA:HB3	53:BW:106:ILE:HD11	1.40	0.99
50:DT:11:GLU:CD	50:DT:11:GLU:H	1.60	0.99
35:BA:586:A:H5'	40:BF:89:VAL:HG21	1.40	0.99
51:BU:92:ARG:HB3	52:BV:11:GLN:NE2	1.76	0.99
1:CA:939:G:H5''	7:CG:102:ARG:HH22	1.24	0.99
1:CA:585:G:H4'	12:CL:8:ASN:HD21	1.28	0.99
35:DA:2348:U:C2'	35:DA:2349:G:H5''	1.92	0.99
52:DV:70:ILE:HB	52:DV:90:PRO:HB2	1.40	0.99
35:BA:1879:C:C2'	35:BA:1880:C:H5''	1.93	0.99
49:BS:13:ARG:H	49:BS:13:ARG:HD2	1.27	0.99
23:CW:3:C:H2'	23:CW:4:G:H5''	1.42	0.99
42:DH:43:VAL:HG11	42:DH:52:VAL:HA	1.44	0.99
56:DZ:24:LEU:HD11	56:DZ:86:VAL:HG23	1.45	0.99
27:B1:32:LYS:HA	35:BA:2396:G:O2'	1.62	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:35:LYS:HG2	38:BD:64:ILE:H	1.28	0.99
43:DI:98:ALA:HB1	43:DI:109:ILE:HB	1.43	0.99
12:AL:83:VAL:HG22	12:AL:84:LEU:H	1.28	0.99
1:CA:1363(A):A:H4'	1:CA:1364:U:H5''	1.41	0.99
49:DS:28:VAL:HG12	49:DS:29:PHE:H	1.25	0.99
26:B0:32:ARG:H	26:B0:35:ASN:HD22	1.05	0.98
35:BA:2125:G:H21	35:BA:2173:A:H62	1.05	0.98
34:D8:32:LEU:C	34:D8:34:TRP:H	1.57	0.98
50:DT:50:ILE:HG23	50:DT:99:LEU:HD12	1.45	0.98
46:BP:30:THR:HG22	46:BP:31:ALA:H	1.27	0.98
52:BV:64:HIS:HB3	52:BV:96:ILE:HA	1.44	0.98
39:DE:24:THR:HG23	39:DE:184:VAL:HG23	1.46	0.98
44:DN:46:VAL:HG13	44:DN:47:ALA:H	1.25	0.98
35:BA:2701:C:H3'	35:BA:2702:U:H5''	1.43	0.98
51:BU:64:ARG:HB3	51:BU:64:ARG:HH21	1.22	0.98
34:D8:46:ARG:NH1	34:D8:46:ARG:HB3	1.78	0.98
35:DA:1494:A:C2'	35:DA:1495:A:H5''	1.93	0.98
35:BA:1590:U:H2'	35:BA:1591:G:H5''	1.42	0.98
35:BA:1657:C:H2'	35:BA:1658:C:C6	1.97	0.98
35:BA:1798:U:H5'	38:BD:259:THR:CG2	1.91	0.98
40:BF:67:GLN:HG3	40:BF:67:GLN:O	1.62	0.98
42:BH:43:VAL:HG11	42:BH:52:VAL:HA	1.41	0.98
19:CS:63:THR:HG22	19:CS:66:MET:HG2	1.45	0.98
35:DA:1586:A:H3'	35:DA:1587:A:C5'	1.94	0.98
34:B8:35:GLN:HA	35:BA:2420:C:OP2	1.64	0.98
10:CJ:34:VAL:HG22	10:CJ:74:ILE:HG22	1.42	0.98
35:DA:2598:A:H5''	38:DD:236:GLY:N	1.76	0.98
35:BA:2036:C:H6	35:BA:2036:C:H5'	1.28	0.98
47:BQ:9:TYR:O	47:BQ:10:ARG:HG3	1.61	0.98
1:CA:1065:U:H5'	1:CA:1190:G:H21	1.29	0.98
3:AC:164:ARG:HB2	3:AC:164:ARG:HH11	1.26	0.98
11:CK:27:ASN:HB2	11:CK:55:LYS:HB3	1.45	0.98
35:DA:27:G:H22	35:DA:512:G:H2'	1.29	0.98
41:DG:37:VAL:HB	41:DG:94:LEU:HB2	1.43	0.98
10:AJ:34:VAL:HG22	10:AJ:74:ILE:HG22	1.44	0.98
39:BE:24:THR:HG23	39:BE:184:VAL:HG23	1.45	0.98
42:BH:18:GLU:HB2	42:BH:25:LYS:HB2	1.42	0.98
13:CM:108:ARG:H	13:CM:108:ARG:HD2	1.28	0.98
35:DA:2415:G:O3'	46:DP:66:GLY:HA3	1.64	0.98
50:BT:40:THR:O	50:BT:41:ARG:HB2	1.63	0.97
2:CB:84:GLU:HB3	2:CB:219:VAL:HG21	1.46	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CP:60:LEU:HD23	16:CP:64:ALA:HB3	1.45	0.97
26:D0:32:ARG:H	26:D0:35:ASN:HD22	1.09	0.97
38:DD:35:LYS:HG2	38:DD:64:ILE:H	1.27	0.97
11:AK:103:LEU:HD22	11:AK:103:LEU:H	1.24	0.97
12:AL:6:THR:HG23	12:AL:9:GLN:H	1.28	0.97
25:AY:150:SER:H	25:AY:153:GLU:HB2	1.29	0.97
45:BO:61:VAL:O	45:BO:84:ALA:HB1	1.62	0.97
35:DA:1798:U:H5'	38:DD:259:THR:CG2	1.92	0.97
35:DA:2600:A:O2'	35:DA:2601:C:H5'	1.64	0.97
54:DX:72:LYS:HE3	54:DX:74:PRO:HB3	1.42	0.97
35:BA:689:A:H2'	35:BA:690:G:H8	1.27	0.97
7:CG:75:VAL:HG12	7:CG:88:PRO:HB3	1.41	0.97
25:CY:129:ILE:HA	25:CY:132:ILE:CD1	1.94	0.97
35:DA:847:U:H2'	35:DA:848:G:H5''	1.46	0.97
36:DB:7:G:H3'	36:DB:8:U:H5''	1.44	0.97
35:DA:389:G:H1	46:DP:71:VAL:HB	1.28	0.97
46:DP:71:VAL:HG13	46:DP:72:PRO:HD3	0.98	0.97
19:AS:63:THR:HG22	19:AS:66:MET:HG2	1.44	0.97
35:BA:1586:A:C3'	35:BA:1587:A:H5''	1.93	0.97
43:BI:98:ALA:HB1	43:BI:109:ILE:HB	1.46	0.97
35:DA:1590:U:H2'	35:DA:1591:G:H5''	1.44	0.97
35:DA:2426:A:H3'	35:DA:2427:C:H5'	1.43	0.97
40:DF:132:VAL:HG22	40:DF:133:ASN:H	1.28	0.97
47:DQ:9:TYR:O	47:DQ:10:ARG:HG3	1.63	0.97
12:AL:37:CYS:HA	12:AL:58:VAL:HG22	1.47	0.97
52:BV:2:PHE:HB3	52:BV:42:GLY:HA2	1.45	0.97
35:DA:2753:A:O2'	35:DA:2754:U:H5'	1.65	0.97
39:DE:116:VAL:HG23	39:DE:120:TRP:HB2	1.47	0.97
35:BA:2645:G:H3'	35:BA:2646:C:H5'	1.45	0.97
8:CH:110:ALA:HB3	8:CH:121:ASP:HB3	1.45	0.97
35:BA:2426:A:H3'	35:BA:2427:C:H5'	1.46	0.97
39:BE:116:VAL:HG23	39:BE:120:TRP:HB2	1.43	0.97
54:BX:60:ARG:HG2	54:BX:74:PRO:HD2	1.46	0.97
35:BA:106:C:H1'	55:BY:2:ARG:NH2	1.80	0.97
2:CB:171:ALA:HA	2:CB:174:VAL:HG23	1.45	0.97
3:CC:164:ARG:HB2	3:CC:164:ARG:HH11	1.24	0.97
20:CT:73:HIS:HB3	20:CT:74:LYS:HD3	1.47	0.97
25:CY:3:LEU:HA	25:CY:6:LEU:HB3	1.41	0.97
35:DA:2645:G:H3'	35:DA:2646:C:H5'	1.44	0.97
39:DE:33:VAL:HG11	39:DE:89:ASP:H	1.27	0.97
41:DG:43:LEU:CD2	41:DG:44:GLY:H	1.77	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:27:ASN:HB2	11:AK:55:LYS:HB3	1.45	0.97
35:BA:2415:G:O3'	46:BP:66:GLY:HA3	1.62	0.97
13:AM:108:ARG:H	13:AM:108:ARG:HD2	1.29	0.97
25:AY:63:PRO:HB2	25:AY:64:ARG:HH12	1.27	0.97
4:CD:145:GLU:HG2	4:CD:184:LYS:HG3	1.46	0.97
47:DQ:55:VAL:HG12	47:DQ:64:ILE:HD12	1.42	0.97
35:BA:2693:A:H2'	35:BA:2694:G:H8	1.30	0.97
12:CL:6:THR:HG23	12:CL:9:GLN:H	1.26	0.97
38:DD:94:LEU:H	38:DD:94:LEU:HD12	1.26	0.97
50:DT:38:ASN:HD22	50:DT:40:THR:H	1.12	0.97
35:BA:925:C:H2'	35:BA:926:A:H5''	1.46	0.96
4:CD:61:LYS:HA	4:CD:203:VAL:HG22	1.45	0.96
56:DZ:149:SER:HB3	56:DZ:173:ALA:HA	1.47	0.96
55:BY:10:GLY:HA2	55:BY:27:VAL:HG13	1.46	0.96
40:DF:103:LYS:HG2	40:DF:106:ARG:HH21	1.28	0.96
44:DN:9:VAL:HG11	44:DN:39:ARG:HH22	1.30	0.96
27:B1:47:GLN:HG2	35:BA:2230:G:C1'	1.94	0.96
44:BN:22:THR:HA	44:BN:61:ARG:HB2	1.47	0.96
11:CK:103:LEU:H	11:CK:103:LEU:HD22	1.27	0.96
35:DA:342:G:H2'	35:DA:343:C:H5''	1.44	0.96
48:BR:48:VAL:HA	48:BR:51:LEU:HD12	1.45	0.96
55:BY:88:LYS:HZ3	55:BY:93:GLY:H	1.01	0.96
25:AY:29:ARG:HE	25:AY:32:ARG:NH2	1.63	0.96
34:B8:46:ARG:HH11	34:B8:46:ARG:CB	1.79	0.96
55:BY:31:LEU:HB3	55:BY:32:PRO:HA	1.47	0.96
56:DZ:125:LEU:HB3	56:DZ:165:VAL:HG22	1.45	0.96
35:BA:27:G:HO2'	35:BA:28:A:H8	1.00	0.96
35:BA:847:U:H2'	35:BA:848:G:H5''	1.46	0.96
50:BT:50:ILE:HG23	50:BT:99:LEU:HD12	1.42	0.96
52:BV:19:LYS:HZ2	52:BV:20:LEU:H	1.09	0.96
27:D1:23:LYS:HB3	27:D1:37:ILE:HD11	1.43	0.96
52:DV:34:GLU:H	52:DV:62:LEU:HB2	1.29	0.96
1:AA:1060:C:H5'	14:AN:45:ARG:HH22	1.30	0.96
44:BN:65:LYS:HA	44:BN:65:LYS:HE2	1.47	0.96
43:DI:91:SER:HB3	43:DI:121:LYS:HE3	1.45	0.96
48:DR:38:VAL:HB	48:DR:39:PRO:HD3	1.48	0.96
50:DT:29:ARG:HB3	50:DT:85:LYS:HA	1.46	0.96
8:AH:110:ALA:HB3	8:AH:121:ASP:HB3	1.45	0.96
48:BR:45:ARG:HG3	48:BR:46:GLY:H	1.30	0.96
41:DG:39:ILE:HA	41:DG:157:ILE:HA	1.45	0.96
1:AA:194:C:H2'	1:AA:195:A:H5''	1.45	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1104:G:H4'	2:AB:111:ARG:NH1	1.81	0.96
5:AE:39:GLY:HA2	5:AE:69:VAL:HB	1.48	0.96
35:BA:1494:A:H2'	35:BA:1495:A:C5'	1.93	0.96
35:BA:2334:G:H5'	49:BS:13:ARG:HG2	1.48	0.96
40:BF:132:VAL:HG22	40:BF:133:ASN:H	1.27	0.96
10:CJ:4:ILE:HB	10:CJ:74:ILE:HD11	1.47	0.96
1:AA:1065:U:H5'	1:AA:1190:G:H21	1.29	0.96
35:DA:925:C:H2'	35:DA:926:A:H5''	1.45	0.96
26:B0:32:ARG:H	26:B0:35:ASN:ND2	1.63	0.95
1:CA:1452:C:H1'	1:CA:1456:G:H22	1.25	0.95
2:CB:112:VAL:O	2:CB:115:LEU:HB3	1.64	0.95
52:DV:2:PHE:HB3	52:DV:42:GLY:HA2	1.48	0.95
27:B1:11:ARG:NH1	27:B1:60:PHE:HA	1.81	0.95
1:CA:1104:G:H4'	2:CB:111:ARG:NH1	1.81	0.95
4:CD:30:LYS:C	4:CD:32:ALA:H	1.68	0.95
49:DS:13:ARG:H	49:DS:13:ARG:HD2	1.27	0.95
52:DV:78:LYS:HD3	52:DV:79:VAL:N	1.81	0.95
35:BA:1024:G:H3'	35:BA:1025:G:H5''	1.48	0.95
45:BO:103:ALA:HA	45:BO:122:LEU:O	1.66	0.95
1:CA:1321:C:C5'	1:CA:1322:C:H5''	1.95	0.95
12:AL:6:THR:O	12:AL:10:LEU:HD23	1.63	0.95
35:BA:342:G:H2'	35:BA:343:C:H5''	1.45	0.95
1:CA:737:A:H2'	1:CA:738:C:C6	2.02	0.95
35:DA:2599:G:OP2	38:DD:236:GLY:HA2	1.65	0.95
35:DA:2562:U:H1'	45:DO:23:ARG:HH12	1.26	0.95
56:DZ:6:LYS:H	56:DZ:6:LYS:HD3	1.31	0.95
38:BD:226:MET:HE2	38:BD:230:ASP:HB3	1.47	0.95
44:BN:46:VAL:HG13	44:BN:47:ALA:H	1.27	0.95
52:BV:34:GLU:H	52:BV:62:LEU:HB2	1.32	0.95
41:DG:46:ALA:HB1	41:DG:88:ILE:HD13	1.49	0.95
35:DA:106:C:H1'	55:DY:2:ARG:NH2	1.81	0.95
5:AE:43:LEU:HD21	5:AE:132:ALA:HB1	1.47	0.95
42:BH:19:VAL:HB	42:BH:44:VAL:HG13	1.47	0.95
34:D8:46:ARG:HH11	34:D8:46:ARG:CB	1.78	0.95
34:B8:46:ARG:NH1	34:B8:46:ARG:HB3	1.79	0.95
5:CE:39:GLY:HA2	5:CE:69:VAL:HB	1.47	0.95
25:CY:84:ARG:HH21	25:CY:92:PRO:HD2	1.32	0.95
1:AA:737:A:H2'	1:AA:738:C:C6	2.00	0.95
36:BB:7:G:H3'	36:BB:8:U:H5''	1.47	0.95
1:CA:383:A:H2'	1:CA:384:G:H5'	1.49	0.95
12:CL:6:THR:O	12:CL:10:LEU:HD23	1.67	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1586:A:C3'	35:DA:1587:A:H5''	1.96	0.95
43:DI:38:LEU:H	43:DI:38:LEU:HD12	1.27	0.95
46:DP:47:ASP:HB3	46:DP:48:PRO:C	1.86	0.95
1:AA:806:C:H2'	1:AA:807:A:H8	1.32	0.95
35:BA:903:C:H2'	35:BA:904:C:C6	2.02	0.95
41:BG:111:LEU:HA	41:BG:114:ILE:HD11	1.46	0.95
43:BI:14:ASP:O	43:BI:17:GLN:HB2	1.67	0.95
50:BT:29:ARG:HB3	50:BT:85:LYS:HA	1.47	0.95
56:BZ:150:LEU:HD22	56:BZ:150:LEU:H	1.27	0.95
1:CA:1349:A:H3'	9:CI:118:LYS:HE3	1.47	0.95
8:AH:23:SER:HA	8:AH:63:LEU:HD23	1.48	0.94
1:AA:522:C:H41	12:AL:53:ARG:HH21	1.07	0.94
39:BE:33:VAL:HG11	39:BE:89:ASP:H	1.32	0.94
52:BV:78:LYS:HD3	52:BV:79:VAL:N	1.82	0.94
11:CK:22:HIS:O	11:CK:28:THR:HG23	1.67	0.94
35:DA:2498:C:O2'	35:DA:2499:C:H5'	1.66	0.94
2:AB:84:GLU:HB3	2:AB:219:VAL:HG21	1.46	0.94
35:BA:2600:A:O2'	35:BA:2601:C:H5'	1.67	0.94
25:CY:125:GLY:O	25:CY:129:ILE:HG13	1.66	0.94
35:DA:2475:C:H42	35:DA:2529:G:H22	1.14	0.94
35:DA:742:G:H2'	35:DA:743:G:H8	1.29	0.94
36:DB:45:A:H8	41:DG:95:ARG:HE	1.12	0.94
46:DP:85:LEU:HD23	46:DP:85:LEU:H	1.29	0.94
10:AJ:4:ILE:HB	10:AJ:74:ILE:HD11	1.46	0.94
18:AR:58:LEU:HD12	18:AR:58:LEU:H	1.28	0.94
35:BA:2740:A:H2'	35:BA:2741:A:C8	2.03	0.94
1:CA:1524:C:H2'	1:CA:1525:G:C8	2.03	0.94
41:DG:43:LEU:HD22	41:DG:44:GLY:H	1.32	0.94
9:AI:28:VAL:HG13	9:AI:64:THR:HA	1.50	0.94
35:BA:2282:G:H1	35:BA:2427:C:H42	1.15	0.94
35:BA:2599:G:OP2	38:BD:236:GLY:HA2	1.67	0.94
50:BT:47:GLY:HA2	50:BT:65:LYS:HB2	1.48	0.94
56:BZ:115:GLY:HA3	56:BZ:175:VAL:O	1.68	0.94
1:CA:559:A:H4'	1:CA:560:U:H5''	1.49	0.94
55:DY:88:LYS:NZ	55:DY:93:GLY:H	1.66	0.94
1:AA:528:C:H41	12:AL:49:ASN:ND2	1.66	0.94
35:BA:2092:U:H4'	35:BA:2093:G:H5''	1.47	0.94
35:BA:742:G:H2'	35:BA:743:G:H8	1.33	0.94
46:BP:39:LYS:CD	46:BP:40:SER:H	1.81	0.94
1:CA:1060:C:H5'	14:CN:45:ARG:HH22	1.32	0.94
2:CB:187:LEU:HD13	2:CB:187:LEU:O	1.67	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1194:U:H2'	1:AA:1195:C:C6	2.02	0.94
35:BA:2538:C:O2'	35:BA:2539:C:H5'	1.68	0.94
38:BD:44:ASN:HB3	38:BD:49:ILE:HA	1.47	0.94
47:BQ:140:ALA:O	56:BZ:72:ARG:HA	1.66	0.94
8:CH:23:SER:HA	8:CH:63:LEU:HD23	1.47	0.94
41:DG:110:ALA:HA	41:DG:140:ILE:HD13	1.48	0.94
45:DO:63:VAL:HG22	45:DO:84:ALA:HA	1.45	0.94
46:DP:47:ASP:HB2	46:DP:51:PHE:HB2	1.48	0.94
55:DY:37:VAL:HG23	55:DY:38:ILE:H	1.32	0.94
35:BA:1494:A:C2'	35:BA:1495:A:H5''	1.96	0.94
35:BA:903:C:H2'	35:BA:904:C:H6	1.32	0.94
31:D5:44:THR:HG21	48:DR:101:ALA:HB2	1.46	0.94
50:DT:40:THR:O	50:DT:41:ARG:HB2	1.64	0.94
13:AM:91:ARG:HH11	19:AS:81:ARG:HH22	1.05	0.94
35:BA:2475:C:H42	35:BA:2529:G:H22	1.13	0.94
54:BX:65:ARG:HE	54:BX:65:ARG:HA	1.33	0.94
1:CA:806:C:H2'	1:CA:807:A:H8	1.32	0.94
35:DA:2125:G:H21	35:DA:2173:A:H62	1.04	0.94
40:DF:53:THR:HG23	40:DF:56:GLU:HB2	1.46	0.94
1:AA:1194:U:H2'	1:AA:1195:C:H6	1.32	0.94
1:CA:878:G:C5'	8:CH:89:PRO:HG2	1.97	0.94
44:DN:22:THR:HA	44:DN:61:ARG:HB2	1.46	0.94
35:BA:2284:C:C2'	35:BA:2285:C:H5''	1.96	0.94
1:AA:1321:C:C5'	1:AA:1322:C:H5''	1.97	0.93
22:AV:30:A:H2'	22:AV:31:U:C6	2.02	0.93
27:B1:37:ILE:H	27:B1:37:ILE:HD12	1.32	0.93
56:DZ:76:LEU:HA	56:DZ:84:GLU:HB2	1.50	0.93
5:AE:147:ASP:HA	5:AE:150:ARG:HH11	1.32	0.93
53:BW:29:LEU:CD2	53:BW:33:ARG:HE	1.81	0.93
12:CL:37:CYS:HA	12:CL:58:VAL:HG22	1.50	0.93
39:BE:78:LEU:H	39:BE:78:LEU:HD23	1.31	0.93
31:B5:44:THR:HG21	48:BR:101:ALA:HB2	1.50	0.93
38:DD:18:VAL:HG23	38:DD:211:ARG:HH21	1.33	0.93
43:DI:14:ASP:O	43:DI:17:GLN:HB2	1.66	0.93
50:DT:106:SER:HA	50:DT:110:ILE:HD13	1.50	0.93
12:AL:55:VAL:HG12	12:AL:56:ALA:H	1.31	0.93
1:CA:1194:U:H2'	1:CA:1195:C:H6	1.33	0.93
27:B1:68:PRO:HG2	27:B1:69:LYS:H	1.29	0.93
55:BY:37:VAL:HG23	55:BY:38:ILE:H	1.33	0.93
1:CA:908:A:H2'	1:CA:909:A:C8	2.04	0.93
1:CA:1397:C:N4	24:CX:22:U:H3'	1.82	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1879:C:C2'	35:DA:1880:C:H5''	1.97	0.93
35:DA:2693:A:H2'	35:DA:2694:G:H8	1.31	0.93
48:DR:45:ARG:HG3	48:DR:46:GLY:H	1.31	0.93
50:DT:28:VAL:HG21	50:DT:47:GLY:N	1.81	0.93
41:BG:63:ILE:HD12	41:BG:64:THR:N	1.83	0.93
27:D1:68:PRO:O	27:D1:70:VAL:N	2.00	0.93
4:AD:145:GLU:HG2	4:AD:184:LYS:HG3	1.48	0.93
42:BH:13:LYS:HA	42:BH:13:LYS:HE2	1.50	0.93
25:CY:147:LEU:HD22	25:CY:149:LEU:HG	1.48	0.93
35:DA:1657:C:H2'	35:DA:1658:C:C6	2.04	0.93
56:BZ:114:GLY:HA3	56:BZ:177:PRO:HB3	1.51	0.93
1:CA:1080:A:H5'	5:CE:14:ARG:HH21	1.30	0.93
25:CY:9:GLU:O	25:CY:12:SER:HB3	1.69	0.93
31:D5:31:VAL:HB	31:D5:32:PRO:HD2	1.50	0.93
34:D8:35:GLN:HA	35:DA:2420:C:OP2	1.68	0.93
35:DA:1902:C:H1'	38:DD:244:ARG:HD3	1.50	0.93
2:AB:172:ILE:HD12	2:AB:172:ILE:H	1.30	0.93
27:B1:47:GLN:HB2	35:BA:397:G:H5''	1.51	0.93
35:DA:1039:G:H1	35:DA:1116:C:H42	1.15	0.93
35:BA:2051:A:O3'	39:BE:141:ILE:HD11	1.69	0.93
19:CS:36:ARG:HH22	19:CS:75:ALA:HB3	1.32	0.93
43:DI:79:ILE:HG12	43:DI:140:LEU:HD11	1.50	0.93
1:AA:625:G:H2'	1:AA:626:U:H6	1.34	0.92
4:AD:30:LYS:C	4:AD:32:ALA:H	1.68	0.92
43:BI:91:SER:HB3	43:BI:121:LYS:HE3	1.52	0.92
7:CG:79:ARG:HE	7:CG:84:ASN:HD21	1.17	0.92
39:DE:78:LEU:HD23	39:DE:78:LEU:H	1.33	0.92
41:DG:39:ILE:HG13	41:DG:157:ILE:HG22	1.51	0.92
4:AD:61:LYS:HA	4:AD:203:VAL:HG22	1.47	0.92
50:BT:106:SER:HA	50:BT:110:ILE:HD13	1.49	0.92
44:DN:65:LYS:HE2	44:DN:65:LYS:HA	1.49	0.92
31:B5:44:THR:HG22	31:B5:45:VAL:H	1.30	0.92
35:BA:2693:A:H2'	35:BA:2694:G:C8	2.05	0.92
55:BY:28:LYS:HZ1	55:BY:37:VAL:HA	1.34	0.92
35:DA:1657:C:H2'	35:DA:1658:C:H6	1.33	0.92
35:DA:2175:C:H2'	35:DA:2176:A:H5''	1.52	0.92
35:DA:2284:C:C2'	35:DA:2285:C:H5''	2.00	0.92
41:DG:111:LEU:HB2	41:DG:112:PRO:CD	1.99	0.92
46:DP:30:THR:HG22	46:DP:31:ALA:H	1.29	0.92
52:DV:38:LEU:HD23	52:DV:39:LEU:N	1.84	0.92
1:AA:559:A:H4'	1:AA:560:U:H5''	1.50	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:AY:63:PRO:HB2	25:AY:64:ARG:NH1	1.82	0.92
44:BN:9:VAL:HG11	44:BN:39:ARG:HH22	1.30	0.92
47:BQ:12:GLN:HG2	47:BQ:73:PRO:HD2	1.51	0.92
27:D1:13:ILE:HG13	27:D1:14:VAL:N	1.83	0.92
35:DA:287:C:N4	35:DA:354:G:H1	1.66	0.92
54:DX:60:ARG:HG2	54:DX:74:PRO:HD2	1.49	0.92
55:DY:87:LYS:HG3	55:DY:89:PHE:H	1.31	0.92
7:AG:79:ARG:HE	7:AG:84:ASN:HD21	1.18	0.92
39:BE:51:PHE:H	39:BE:74:PRO:CG	1.82	0.92
46:BP:47:ASP:HB2	46:BP:51:PHE:HB2	1.50	0.92
2:CB:71:VAL:HG22	2:CB:93:VAL:N	1.83	0.92
12:CL:69:TYR:O	12:CL:71:PRO:HD3	1.69	0.92
27:D1:13:ILE:HG13	27:D1:14:VAL:H	1.35	0.92
35:DA:1747(A):G:C2'	35:DA:1748:G:H5''	2.00	0.92
1:AA:452:A:HO2'	1:AA:453:A:H8	0.95	0.92
1:AA:1349:A:H3'	9:AI:118:LYS:HE3	1.50	0.92
35:BA:2314:C:H5'	41:BG:38:VAL:HG11	1.52	0.92
35:BA:271(D):G:H1	35:BA:271(T):C:H42	0.94	0.92
53:BW:29:LEU:HD21	53:BW:33:ARG:NE	1.83	0.92
5:CE:98:THR:HB	5:CE:117:ASP:HB3	1.52	0.92
18:CR:58:LEU:H	18:CR:58:LEU:HD12	1.33	0.92
31:D5:44:THR:HG22	31:D5:45:VAL:H	1.32	0.92
27:B1:64:ALA:O	27:B1:67:ILE:HG13	1.70	0.92
29:B3:56:VAL:HG12	29:B3:57:GLU:H	1.35	0.92
34:B8:30:ARG:HH21	46:BP:62:LEU:HB2	1.35	0.92
25:CY:126:ARG:HA	25:CY:129:ILE:HD12	1.52	0.92
35:DA:1024:G:H3'	35:DA:1025:G:H5''	1.48	0.92
31:B5:31:VAL:HB	31:B5:32:PRO:HD2	1.51	0.92
35:BA:287:C:N4	35:BA:354:G:H1	1.68	0.92
35:BA:364:C:H2'	35:BA:365:C:H5''	1.52	0.92
41:BG:46:ALA:HB1	41:BG:88:ILE:HD13	1.50	0.92
1:CA:522:C:H41	12:CL:53:ARG:HH21	1.08	0.92
54:DX:25:LYS:NZ	54:DX:87:GLN:H	1.67	0.92
48:BR:38:VAL:HB	48:BR:39:PRO:HD3	1.48	0.92
53:BW:78:GLU:OE2	53:BW:99:ARG:HD2	1.69	0.92
56:BZ:56:VAL:HG22	56:BZ:70:LEU:HG	1.52	0.92
35:DA:2538:C:O2'	35:DA:2539:C:H5'	1.70	0.92
35:DA:2334:G:H5'	49:DS:13:ARG:HG2	1.50	0.92
4:AD:18:LYS:HZ3	4:AD:33:MET:HB3	1.34	0.92
51:BU:66:ASN:ND2	51:BU:70:ARG:HH21	1.66	0.92
56:BZ:149:SER:HB3	56:BZ:173:ALA:HA	1.50	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:172:ILE:H	2:CB:172:ILE:HD12	1.32	0.92
35:DA:2758:A:C2'	35:DA:2759:G:H5''	1.99	0.92
56:DZ:110:GLY:O	56:DZ:112:ARG:N	2.02	0.92
56:DZ:166:SER:HB2	56:DZ:167:PRO:CA	1.99	0.92
1:AA:1347:G:N2	1:AA:1373:G:H2'	1.86	0.91
35:BA:1791:A:H5'	38:BD:206:LEU:HD13	1.51	0.91
46:BP:85:LEU:HD23	46:BP:85:LEU:H	1.32	0.91
51:BU:106:PHE:HA	51:BU:109:LEU:HD12	1.52	0.91
55:BY:88:LYS:NZ	55:BY:93:GLY:H	1.68	0.91
27:D1:25:LYS:HG3	27:D1:37:ILE:HG21	1.51	0.91
35:DA:1348:G:C2'	35:DA:1349:A:H5''	2.00	0.91
41:DG:91:ARG:HD2	41:DG:92:VAL:N	1.85	0.91
50:DT:47:GLY:HA2	50:DT:65:LYS:HB2	1.51	0.91
51:DU:92:ARG:HG2	51:DU:95:LEU:H	1.35	0.91
35:BA:2753:A:O2'	35:BA:2754:U:H5'	1.69	0.91
47:DQ:108:GLY:HA3	56:DZ:116:VAL:HG21	1.49	0.91
4:AD:108:LEU:HB3	4:AD:110:PHE:HE1	1.33	0.91
26:B0:70:GLN:HG2	26:B0:71:ASP:N	1.85	0.91
38:BD:35:LYS:HG2	38:BD:64:ILE:N	1.84	0.91
43:BI:92:VAL:HG11	43:BI:120:ILE:HD12	1.51	0.91
2:CB:236:TYR:HA	2:CB:239:VAL:CG2	2.01	0.91
11:CK:29:ILE:HG22	11:CK:44:SER:HB3	1.51	0.91
35:DA:1177:A:H3'	35:DA:1177:A:OP2	1.70	0.91
53:DW:29:LEU:CD2	53:DW:33:ARG:HE	1.83	0.91
28:D2:26:ARG:HH22	54:DX:6:ASP:HA	1.35	0.91
1:AA:1169:A:H2'	1:AA:1170:A:C8	2.04	0.91
27:B1:51:VAL:O	27:B1:60:PHE:HB2	1.71	0.91
35:BA:1039:G:H1	35:BA:1116:C:H42	1.15	0.91
44:BN:65:LYS:O	44:BN:69:GLN:HB2	1.71	0.91
9:CI:28:VAL:HG13	9:CI:64:THR:HA	1.48	0.91
26:D0:32:ARG:H	26:D0:35:ASN:ND2	1.67	0.91
35:DA:364:C:H2'	35:DA:365:C:H5''	1.51	0.91
53:DW:29:LEU:HD21	53:DW:33:ARG:NE	1.85	0.91
1:CA:1169:A:H2'	1:CA:1170:A:C8	2.04	0.91
1:CA:1194:U:H2'	1:CA:1195:C:C6	2.05	0.91
1:CA:1403:C:H1'	1:CA:1500:A:N1	1.85	0.91
41:DG:11:TYR:O	41:DG:16:ARG:HB2	1.69	0.91
46:BP:85:LEU:HA	46:BP:88:LEU:HB3	1.53	0.91
35:BA:2820:A:H4'	48:BR:5:LYS:HE2	1.50	0.91
1:CA:954:G:H4'	13:CM:120:LYS:HG3	1.53	0.91
33:D7:24:THR:HG23	33:D7:27:GLY:HA3	1.50	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:44:ASN:HB3	38:DD:49:ILE:HA	1.52	0.91
44:DN:65:LYS:O	44:DN:69:GLN:HB2	1.71	0.91
45:DO:103:ALA:HA	45:DO:122:LEU:O	1.69	0.91
54:DX:65:ARG:HA	54:DX:65:ARG:HE	1.35	0.91
2:AB:71:VAL:HG22	2:AB:93:VAL:N	1.84	0.91
35:BA:2297:C:H2'	35:BA:2298:A:H5'	1.53	0.91
47:BQ:39:PRO:CB	47:BQ:99:PRO:HD3	2.01	0.91
52:BV:62:LEU:HD22	52:BV:96:ILE:HD13	1.53	0.91
55:BY:28:LYS:NZ	55:BY:30:VAL:HA	1.86	0.91
1:CA:1347:G:N2	1:CA:1373:G:H2'	1.86	0.91
38:DD:161:THR:HG1	38:DD:196:VAL:HG21	1.31	0.91
40:DF:158:THR:HG21	40:DF:163:VAL:HB	1.53	0.91
12:AL:69:TYR:O	12:AL:71:PRO:HD3	1.70	0.91
16:AP:4:ILE:HB	16:AP:66:PRO:HB3	1.51	0.91
35:BA:27:G:H22	35:BA:512:G:H2'	1.34	0.91
40:BF:114:VAL:HG23	40:BF:115:ALA:H	1.36	0.91
42:BH:43:VAL:HG21	42:BH:52:VAL:HG22	1.52	0.91
47:BQ:141:GLN:HB3	56:BZ:70:LEU:HD13	1.53	0.91
2:CB:47:THR:HG23	2:CB:202:PRO:HG2	1.52	0.91
35:DA:144:C:H2'	35:DA:145:G:H8	1.35	0.91
35:DA:272(D):G:H1	35:DA:364:C:H42	1.19	0.91
35:DA:549:G:H2'	35:DA:551:G:H5''	1.50	0.91
1:AA:954:G:H4'	13:AM:120:LYS:HG3	1.51	0.91
19:AS:53:ASN:HD21	19:AS:56:GLN:H	1.18	0.91
35:BA:1986:A:H3'	35:BA:1987:G:H5''	1.53	0.91
11:CK:22:HIS:HB3	11:CK:29:ILE:HG13	1.52	0.91
46:DP:115:LEU:HA	46:DP:134:ALA:HB1	1.53	0.91
8:AH:12:ARG:NH1	8:AH:27:PRO:HD3	1.85	0.91
35:BA:1177:A:OP2	35:BA:1177:A:H3'	1.70	0.91
35:BA:26:G:H1'	35:BA:515:A:H61	1.35	0.91
40:BF:103:LYS:HG2	40:BF:106:ARG:HH21	1.35	0.91
50:BT:38:ASN:HD22	50:BT:40:THR:H	1.13	0.91
5:CE:147:ASP:HA	5:CE:150:ARG:HH11	1.36	0.91
35:DA:2051:A:O3'	39:DE:141:ILE:HD11	1.71	0.91
2:AB:71:VAL:HG13	2:AB:93:VAL:O	1.71	0.90
18:CR:56:THR:HB	18:CR:58:LEU:HD13	1.53	0.90
35:DA:740:U:H2'	35:DA:741:G:C8	2.05	0.90
46:DP:39:LYS:CD	46:DP:40:SER:H	1.85	0.90
35:BA:93:G:H2'	35:BA:94:C:C6	2.07	0.90
54:BX:55:ASN:C	54:BX:77:LYS:HG3	1.92	0.90
4:CD:108:LEU:HB3	4:CD:110:PHE:HE1	1.34	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:18:LYS:HZ3	4:CD:33:MET:HB3	1.34	0.90
38:DD:36:PRO:HG3	38:DD:61:LEU:HD21	1.52	0.90
1:AA:1116:C:H2'	1:AA:1117:G:H5''	1.52	0.90
20:AT:73:HIS:HB3	20:AT:74:LYS:HD3	1.51	0.90
34:B8:22:VAL:HB	34:B8:53:PRO:HB2	1.53	0.90
38:BD:14:ARG:HB2	38:BD:14:ARG:HH11	1.36	0.90
38:BD:76:PRO:HG2	38:BD:98:VAL:HG21	1.52	0.90
52:BV:38:LEU:HD23	52:BV:39:LEU:N	1.87	0.90
55:BY:87:LYS:HG3	55:BY:89:PHE:H	1.32	0.90
5:CE:43:LEU:HD21	5:CE:132:ALA:HB1	1.50	0.90
8:CH:33:GLU:HA	8:CH:36:LEU:HD12	1.52	0.90
41:DG:114:ILE:CG2	41:DG:117:PHE:HB2	2.02	0.90
42:DH:13:LYS:HA	42:DH:13:LYS:HE2	1.51	0.90
1:AA:908:A:H2'	1:AA:909:A:C8	2.06	0.90
35:BA:549:G:H2'	35:BA:551:G:H5''	1.50	0.90
44:BN:74:ARG:NH2	44:BN:101:HIS:HB3	1.87	0.90
44:BN:120:LEU:HD11	44:BN:122:VAL:HG23	1.54	0.90
52:BV:32:THR:HG22	52:BV:33:VAL:H	1.37	0.90
34:D8:22:VAL:HB	34:D8:53:PRO:HB2	1.52	0.90
35:DA:903:C:H2'	35:DA:904:C:C6	2.05	0.90
38:DD:35:LYS:HG2	38:DD:64:ILE:N	1.86	0.90
50:DT:23:ARG:O	50:DT:25:GLY:N	2.04	0.90
35:BA:2426:A:H3'	35:BA:2427:C:C5'	2.02	0.90
52:BV:18:LEU:HD22	52:BV:19:LYS:N	1.86	0.90
31:D5:20:ARG:HH12	53:DW:15:ARG:NH2	1.68	0.90
41:DG:7:LEU:HA	41:DG:10:LYS:HB2	1.51	0.90
51:DU:17:ILE:HA	51:DU:20:LEU:HD23	1.51	0.90
3:AC:15:THR:HG22	3:AC:16:ARG:NH1	1.86	0.90
35:BA:2758:A:C2'	35:BA:2759:G:H5''	2.01	0.90
55:BY:31:LEU:HB3	55:BY:32:PRO:CA	2.02	0.90
56:BZ:110:GLY:C	56:BZ:112:ARG:H	1.69	0.90
15:CO:9:GLN:HB3	15:CO:13:GLN:HE21	1.35	0.90
45:DO:61:VAL:O	45:DO:84:ALA:HB1	1.70	0.90
51:DU:64:ARG:CB	51:DU:64:ARG:HH21	1.83	0.90
55:DY:45:VAL:HG22	55:DY:62:GLU:HB2	1.54	0.90
1:AA:1469:G:H2'	1:AA:1470:G:C8	2.05	0.90
1:AA:266:G:H5''	1:AA:268:C:H41	1.36	0.90
1:AA:383:A:H2'	1:AA:384:G:H5'	1.52	0.90
13:AM:118:ALA:HB1	13:AM:119:GLY:N	1.86	0.90
2:CB:71:VAL:HG13	2:CB:93:VAL:O	1.72	0.90
43:DI:144:VAL:HG12	43:DI:145:VAL:HG23	1.54	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:34:LEU:HD11	47:DQ:129:THR:HB	1.53	0.90
49:DS:25:ARG:HH21	49:DS:89:ARG:HH12	1.17	0.90
1:AA:489:C:H2'	1:AA:490:G:H8	1.36	0.90
2:AB:187:LEU:HD13	2:AB:187:LEU:O	1.72	0.90
2:AB:236:TYR:HA	2:AB:239:VAL:CG2	2.02	0.90
35:BA:1747(A):G:C2'	35:BA:1748:G:H5''	2.02	0.90
46:BP:101:VAL:HG23	46:BP:107:LYS:HA	1.54	0.90
50:BT:28:VAL:HG21	50:BT:47:GLY:N	1.86	0.90
54:BX:73:ARG:H	54:BX:74:PRO:CD	1.85	0.90
8:CH:129:VAL:HG23	8:CH:130:GLY:H	1.37	0.90
8:CH:82:HIS:HD2	8:CH:138:TRP:HE1	1.18	0.90
27:D1:75:GLU:HB2	27:D1:76:ARG:NH2	1.87	0.90
34:D8:53:PRO:HA	34:D8:56:GLU:OE1	1.71	0.90
35:DA:2092:U:H4'	35:DA:2093:G:H5''	1.51	0.90
41:DG:114:ILE:HG22	41:DG:115:ARG:H	1.35	0.90
47:DQ:133:ARG:O	47:DQ:134:ARG:HB2	1.68	0.90
52:DV:62:LEU:HD22	52:DV:96:ILE:HD13	1.53	0.90
4:AD:11:LEU:C	4:AD:13:ARG:H	1.74	0.90
19:AS:70:LYS:HZ3	19:AS:70:LYS:HB3	1.36	0.90
1:CA:452:A:HO2'	1:CA:453:A:H8	0.95	0.90
35:DA:2426:A:H3'	35:DA:2427:C:C5'	2.01	0.90
51:DU:66:ASN:ND2	51:DU:70:ARG:HH21	1.69	0.90
55:DY:28:LYS:NZ	55:DY:30:VAL:HA	1.86	0.90
55:DY:31:LEU:HB3	55:DY:32:PRO:HA	1.50	0.90
56:DZ:27:VAL:HA	56:DZ:36:LYS:HA	1.53	0.90
1:AA:955:U:H1'	1:AA:1227:A:H61	1.34	0.90
11:AK:22:HIS:HB3	11:AK:29:ILE:HG13	1.52	0.90
40:BF:63:LYS:NZ	40:BF:67:GLN:HB3	1.87	0.90
46:BP:115:LEU:HA	46:BP:134:ALA:HB1	1.51	0.90
50:BT:120:ARG:O	50:BT:124:ASP:HB2	1.71	0.90
7:CG:69:VAL:HA	7:CG:138:LYS:HD2	1.55	0.90
8:CH:30:ARG:HH11	8:CH:30:ARG:HB3	1.35	0.90
35:DA:2310:A:O2'	35:DA:2311:A:H5'	1.72	0.90
46:DP:75:ILE:HD12	46:DP:75:ILE:H	1.36	0.90
8:AH:30:ARG:HB3	8:AH:30:ARG:HH11	1.34	0.89
23:AW:10:G:N2	23:AW:27:G:H1'	1.86	0.89
35:BA:1403:C:H5''	35:BA:1471:A:H1'	1.53	0.89
35:BA:2850:A:H2'	35:BA:2851:A:H8	1.35	0.89
35:BA:287:C:H42	35:BA:354:G:H1	0.90	0.89
40:BF:129:PHE:HA	40:BF:142:TRP:HE1	1.36	0.89
50:BT:13:ARG:HA	50:BT:13:ARG:NE	1.87	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:979:C:H3'	1:CA:980:C:C5'	2.01	0.89
7:CG:60:LYS:HA	7:CG:63:LYS:HB3	1.54	0.89
9:CI:50:LEU:HD21	9:CI:81:ILE:HG21	1.54	0.89
16:CP:20:VAL:CG2	16:CP:32:TYR:HB2	2.02	0.89
35:DA:491:G:H2'	35:DA:492:A:H8	1.34	0.89
47:DQ:12:GLN:HG2	47:DQ:73:PRO:HD2	1.52	0.89
49:DS:89:ARG:HA	49:DS:89:ARG:HE	1.37	0.89
15:AO:9:GLN:HB3	15:AO:13:GLN:HE21	1.36	0.89
18:AR:56:THR:HB	18:AR:58:LEU:HD13	1.52	0.89
54:BX:25:LYS:NZ	54:BX:87:GLN:H	1.69	0.89
38:DD:131:LEU:HA	38:DD:190:TYR:CE2	2.07	0.89
38:DD:35:LYS:HB3	38:DD:63:ARG:HA	1.54	0.89
42:DH:66:GLY:HA2	42:DH:69:ARG:HB2	1.54	0.89
51:DU:66:ASN:HD21	51:DU:70:ARG:HH21	0.90	0.89
5:AE:33:VAL:HG22	5:AE:43:LEU:HD13	1.55	0.89
9:AI:50:LEU:HD21	9:AI:81:ILE:HG21	1.53	0.89
33:B7:24:THR:HG23	33:B7:27:GLY:HA3	1.53	0.89
35:BA:1348:G:C2'	35:BA:1349:A:H5''	2.01	0.89
35:BA:2591:C:H2'	35:BA:2592:G:C8	2.07	0.89
38:BD:35:LYS:HB3	38:BD:63:ARG:HA	1.55	0.89
56:BZ:17:ALA:HA	56:BZ:20:ARG:HB3	1.51	0.89
35:DA:1019:U:H3	35:DA:1142(A):A:H62	1.17	0.89
39:DE:116:VAL:HG21	39:DE:122:PHE:CD2	2.08	0.89
44:DN:120:LEU:HD11	44:DN:122:VAL:HG23	1.52	0.89
46:DP:112:LEU:HD11	46:DP:114:ILE:HG23	1.54	0.89
50:DT:80:SER:HB3	50:DT:81:PRO:HD2	1.52	0.89
1:AA:979:C:H3'	1:AA:980:C:C5'	2.01	0.89
8:AH:97:VAL:HG13	8:AH:98:LYS:H	1.37	0.89
35:BA:2712:U:O2'	35:BA:2712(A):A:H5''	1.72	0.89
40:BF:158:THR:HG21	40:BF:163:VAL:HB	1.52	0.89
42:BH:123:PHE:HA	42:BH:133:VAL:HG22	1.53	0.89
49:BS:25:ARG:HH21	49:BS:89:ARG:HH12	1.20	0.89
1:CA:955:U:H1'	1:CA:1227:A:H61	1.35	0.89
27:D1:9:GLY:H	27:D1:48:LYS:HZ3	1.17	0.89
35:DA:2693:A:H2'	35:DA:2694:G:C8	2.07	0.89
53:DW:78:GLU:OE2	53:DW:99:ARG:HD2	1.73	0.89
27:B1:58:ILE:HD13	27:B1:59:THR:N	1.86	0.89
19:CS:53:ASN:HD21	19:CS:56:GLN:H	1.20	0.89
1:AA:678:U:H2'	1:AA:679:C:C6	2.07	0.89
27:B1:48:LYS:HG3	27:B1:49:VAL:H	1.37	0.89
43:BI:144:VAL:HG12	43:BI:145:VAL:HG23	1.53	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BI:79:ILE:HG12	43:BI:140:LEU:HD11	1.54	0.89
2:CB:96:ARG:N	2:CB:96:ARG:HD2	1.88	0.89
27:D1:21:ARG:HD3	27:D1:21:ARG:C	1.93	0.89
35:DA:2740:A:H2'	35:DA:2741:A:C8	2.07	0.89
35:DA:689:A:H2'	35:DA:690:G:C8	2.08	0.89
50:DT:120:ARG:O	50:DT:124:ASP:HB2	1.73	0.89
56:DZ:97:GLU:HB3	56:DZ:125:LEU:HD21	1.54	0.89
1:AA:707:C:H4'	11:AK:20:TYR:CD1	2.06	0.89
40:BF:53:THR:H	40:BF:56:GLU:HB2	1.34	0.89
46:BP:77:ARG:HB2	46:BP:78:PRO:HD2	1.54	0.89
49:BS:89:ARG:HA	49:BS:89:ARG:HE	1.36	0.89
54:BX:60:ARG:HG3	54:BX:72:LYS:H	1.35	0.89
9:CI:19:LEU:HD23	9:CI:61:ALA:HA	1.55	0.89
35:DA:2276:G:H5'	47:DQ:84:GLY:HA2	1.54	0.89
35:DA:271(D):G:H1	35:DA:271(T):C:H42	0.96	0.89
55:DY:88:LYS:HZ3	55:DY:93:GLY:H	0.92	0.89
8:AH:20:TYR:HA	8:AH:65:TYR:CE2	2.08	0.89
38:BD:131:LEU:HD13	38:BD:136:ILE:HG12	1.54	0.89
56:BZ:103:ARG:HB2	56:BZ:138:GLU:HA	1.51	0.89
1:CA:528:C:H41	12:CL:49:ASN:ND2	1.71	0.89
35:DA:2712:U:O2'	35:DA:2712(A):A:H5''	1.72	0.89
35:DA:903:C:H2'	35:DA:904:C:H6	1.36	0.89
40:DF:39:TRP:CD1	40:DF:101:LEU:HB2	2.08	0.89
2:AB:165:VAL:HG23	2:AB:166:ASP:N	1.86	0.89
56:BZ:58:VAL:HG22	56:BZ:68:PRO:HA	1.55	0.89
1:CA:707:C:H4'	11:CK:20:TYR:CD1	2.08	0.89
4:CD:11:LEU:C	4:CD:13:ARG:H	1.72	0.89
8:CH:12:ARG:NH1	8:CH:27:PRO:HD3	1.87	0.89
35:DA:26:G:H1'	35:DA:515:A:H61	1.38	0.89
2:AB:42:ILE:HD11	2:AB:202:PRO:HB2	1.55	0.89
35:BA:2175:C:H2'	35:BA:2176:A:H5''	1.53	0.89
27:B1:47:GLN:CG	35:BA:2230:G:H1'	1.96	0.89
38:BD:44:ASN:HB2	38:BD:48:ARG:O	1.73	0.89
47:BQ:34:LEU:HD11	47:BQ:129:THR:HB	1.53	0.89
1:CA:932:C:H5''	7:CG:3:ARG:HD2	1.53	0.89
35:DA:287:C:H42	35:DA:354:G:H1	0.89	0.89
35:BA:2672:G:H2'	35:BA:2673:G:H5''	1.55	0.88
41:BG:72:ARG:HB3	41:BG:86:MET:H	1.37	0.88
52:BV:72:VAL:HG12	52:BV:73:SER:H	1.38	0.88
41:DG:37:VAL:HG23	41:DG:99:MET:HG3	1.55	0.88
55:DY:45:VAL:HG13	55:DY:62:GLU:OE2	1.71	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:100:ARG:HH12	4:AD:137:SER:HA	1.38	0.88
5:AE:98:THR:HB	5:AE:117:ASP:HB3	1.53	0.88
35:DA:2297:C:H2'	35:DA:2298:A:H5'	1.52	0.88
42:DH:123:PHE:HA	42:DH:133:VAL:HG22	1.56	0.88
50:DT:65:LYS:HA	50:DT:65:LYS:NZ	1.87	0.88
51:DU:106:PHE:HA	51:DU:109:LEU:HD12	1.52	0.88
2:AB:96:ARG:N	2:AB:96:ARG:HD2	1.87	0.88
41:BG:153:ARG:HG3	41:BG:153:ARG:HH11	1.37	0.88
50:BT:80:SER:HB3	50:BT:81:PRO:HD2	1.52	0.88
27:D1:86:SER:HA	27:D1:89:GLU:OE1	1.73	0.88
35:DA:1255:U:C5'	35:DA:1256:G:H5''	2.03	0.88
38:DD:44:ASN:HB2	38:DD:48:ARG:O	1.72	0.88
42:DH:43:VAL:HG21	42:DH:52:VAL:HG22	1.55	0.88
8:AH:33:GLU:HA	8:AH:36:LEU:HD12	1.54	0.88
19:AS:16:LEU:O	19:AS:20:LEU:HG	1.73	0.88
31:B5:20:ARG:HH12	53:BW:15:ARG:NH2	1.71	0.88
34:B8:52:LYS:N	34:B8:53:PRO:HD2	1.87	0.88
35:BA:1986:A:C3'	35:BA:1987:G:H5''	2.02	0.88
38:BD:18:VAL:HG23	38:BD:211:ARG:HH21	1.38	0.88
38:BD:94:LEU:HD12	38:BD:94:LEU:H	1.36	0.88
39:BE:51:PHE:N	39:BE:74:PRO:HG3	1.87	0.88
43:BI:2:LYS:HB2	43:BI:39:ALA:HB2	1.56	0.88
46:BP:23:PRO:HB2	46:BP:33:ARG:CD	2.03	0.88
1:CA:625:G:H2'	1:CA:626:U:H6	1.38	0.88
39:DE:137:HIS:HB3	39:DE:138:PRO:HD2	1.54	0.88
52:DV:72:VAL:HG12	52:DV:73:SER:H	1.36	0.88
34:B8:53:PRO:HA	34:B8:56:GLU:OE1	1.74	0.88
41:BG:27:ASN:ND2	41:BG:29:TRP:HB2	1.89	0.88
55:BY:45:VAL:HG22	55:BY:62:GLU:HB2	1.56	0.88
8:CH:97:VAL:HG13	8:CH:98:LYS:H	1.39	0.88
19:CS:16:LEU:O	19:CS:20:LEU:HG	1.73	0.88
27:D1:67:ILE:N	27:D1:68:PRO:HD2	1.89	0.88
35:DA:570:G:H2'	35:DA:2030:A:N7	1.89	0.88
41:DG:32:PRO:HB2	41:DG:172:LEU:HD13	1.54	0.88
14:AN:8:GLU:HB2	14:AN:12:ARG:HH11	1.37	0.88
35:BA:144:C:H2'	35:BA:145:G:H8	1.37	0.88
38:BD:36:PRO:HG3	38:BD:61:LEU:HD21	1.54	0.88
42:BH:19:VAL:HG21	42:BH:44:VAL:HA	1.55	0.88
6:CF:45:LEU:HD12	6:CF:46:ARG:H	1.36	0.88
35:DA:521:G:H2'	35:DA:522:G:H8	1.38	0.88
46:DP:7:ARG:HB3	46:DP:8:PRO:HD3	1.54	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:13:ARG:HA	50:DT:13:ARG:NE	1.89	0.88
25:AY:70:SER:HB3	25:AY:76:LEU:HD12	1.55	0.88
35:BA:1930:G:N2	35:BA:1968:G:H2'	1.88	0.88
47:BQ:133:ARG:O	47:BQ:134:ARG:HB2	1.73	0.88
12:CL:6:THR:HG22	12:CL:9:GLN:HG3	1.54	0.88
28:D2:60:LEU:HG	28:D2:61:LEU:H	1.38	0.88
35:DA:1230:C:H2'	35:DA:1231:G:H8	1.37	0.88
35:DA:1403:C:H5''	35:DA:1471:A:H1'	1.54	0.88
35:DA:2282:G:H1	35:DA:2427:C:H42	1.21	0.88
51:DU:92:ARG:HD3	51:DU:94:ASN:HD22	1.38	0.88
1:AA:591:U:H2'	1:AA:592:G:C8	2.09	0.88
6:AF:68:PRO:HG3	6:AF:71:ARG:HH21	1.37	0.88
23:AW:73:A:OP1	35:BA:1852:C:H5'	1.74	0.88
3:AC:79:ARG:HH12	11:CK:99:GLN:HB3	1.38	0.88
35:DA:708:C:H42	35:DA:723:G:H1	1.21	0.88
40:DF:53:THR:H	40:DF:56:GLU:HB2	1.38	0.88
56:DZ:48:PHE:HE2	56:DZ:71:VAL:HG21	1.36	0.88
3:AC:105:GLU:HG2	3:AC:106:VAL:H	1.38	0.88
25:AY:3:LEU:H	25:AY:3:LEU:HD12	1.39	0.88
52:BV:71:LEU:HD13	52:BV:72:VAL:H	1.38	0.88
1:CA:673:G:H2'	1:CA:674:G:C8	2.08	0.88
26:D0:70:GLN:HG2	26:D0:71:ASP:N	1.85	0.88
35:DA:93:G:H2'	35:DA:94:C:C6	2.08	0.88
39:DE:51:PHE:H	39:DE:74:PRO:CG	1.85	0.88
35:BA:570:G:H2'	35:BA:2030:A:N7	1.86	0.88
46:BP:127:ALA:HB3	46:BP:130:PHE:CE2	2.08	0.88
50:BT:65:LYS:HA	50:BT:65:LYS:NZ	1.89	0.88
1:CA:1458:G:H2'	1:CA:1459:C:C6	2.09	0.88
42:DH:19:VAL:HG21	42:DH:44:VAL:HA	1.53	0.88
4:AD:100:ARG:NH1	4:AD:137:SER:HA	1.89	0.87
8:AH:6:ILE:HG21	8:AH:85:ARG:NH1	1.89	0.87
1:AA:878:G:C5'	8:AH:89:PRO:HG2	2.03	0.87
35:BA:2310:A:O2'	35:BA:2311:A:H5'	1.73	0.87
42:BH:66:GLY:HA2	42:BH:69:ARG:HB2	1.54	0.87
43:BI:72:LEU:HD12	43:BI:138:ILE:CD1	2.04	0.87
23:CW:1:C:H42	23:CW:74:A:H61	1.20	0.87
25:CY:18:LEU:HG	25:CY:19:GLU:N	1.89	0.87
35:DA:1986:A:C3'	35:DA:1987:G:H5''	2.04	0.87
35:DA:484:C:H2'	35:DA:485:C:C6	2.08	0.87
39:DE:52:LEU:HB2	39:DE:76:ARG:HB2	1.55	0.87
43:BI:102:SER:HB2	43:BI:109:ILE:HG12	1.57	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2672:G:H2'	35:DA:2673:G:H5''	1.54	0.87
56:DZ:26:GLY:HA3	56:DZ:37:VAL:O	1.74	0.87
16:AP:20:VAL:CG2	16:AP:32:TYR:HB2	2.05	0.87
46:BP:112:LEU:HD11	46:BP:114:ILE:HG23	1.57	0.87
25:CY:15:GLN:HA	25:CY:18:LEU:HD23	1.57	0.87
27:D1:9:GLY:H	27:D1:48:LYS:NZ	1.71	0.87
43:DI:2:LYS:HB2	43:DI:39:ALA:HB2	1.55	0.87
44:DN:89:LYS:O	44:DN:93:THR:HG22	1.74	0.87
46:DP:85:LEU:HA	46:DP:88:LEU:HB3	1.56	0.87
56:DZ:149:SER:CB	56:DZ:173:ALA:HA	2.04	0.87
35:BA:588:U:H2'	35:BA:589:C:C6	2.10	0.87
40:BF:178:PRO:HG2	40:BF:179:GLU:H	1.39	0.87
41:BG:178:PHE:HB3	41:BG:180:PHE:HE1	1.39	0.87
42:BH:70:THR:HG22	42:BH:74:ASN:ND2	1.90	0.87
55:BY:75:ILE:HD13	55:BY:76:CYS:H	1.38	0.87
1:CA:1169:A:H2'	1:CA:1170:A:H8	1.38	0.87
9:CI:114:TYR:H	9:CI:114:TYR:HD2	1.20	0.87
35:DA:1577:C:H2'	35:DA:1578:U:C6	2.09	0.87
35:DA:769:G:O2'	35:DA:770:G:H5'	1.75	0.87
46:DP:101:VAL:HG23	46:DP:107:LYS:HA	1.54	0.87
56:DZ:56:VAL:HA	56:DZ:70:LEU:CD2	1.99	0.87
39:BE:30:PRO:HD3	39:BE:180:ASN:HD21	1.39	0.87
1:CA:1412:C:H2'	1:CA:1413:A:C8	2.09	0.87
38:DD:14:ARG:HH11	38:DD:14:ARG:HB2	1.37	0.87
41:DG:130:ASN:ND2	41:DG:161:THR:H	1.72	0.87
43:DI:72:LEU:HD12	43:DI:138:ILE:CD1	2.04	0.87
54:DX:60:ARG:HG3	54:DX:72:LYS:H	1.38	0.87
55:DY:75:ILE:HD13	55:DY:76:CYS:H	1.38	0.87
12:AL:6:THR:HG22	12:AL:9:GLN:HG3	1.55	0.87
28:B2:50:ILE:HG23	28:B2:54:LYS:HD3	1.57	0.87
35:BA:2831:G:H1'	35:BA:2883:A:H2'	1.57	0.87
28:B2:53:LEU:HD12	35:BA:76:C:O3'	1.74	0.87
45:BO:101:PRO:O	45:BO:102:VAL:HG13	1.75	0.87
40:DF:178:PRO:HG2	40:DF:179:GLU:H	1.37	0.87
43:DI:91:SER:CB	43:DI:119:PRO:HB2	2.04	0.87
16:AP:49:LEU:HD11	16:AP:51:VAL:HG23	1.56	0.87
17:AQ:97:SER:O	17:AQ:98:LEU:HG	1.75	0.87
35:BA:521:G:H2'	35:BA:522:G:H8	1.40	0.87
51:BU:92:ARG:HG2	51:BU:95:LEU:H	1.40	0.87
16:CP:4:ILE:HB	16:CP:66:PRO:HB3	1.54	0.87
28:D2:14:ARG:HE	28:D2:14:ARG:C	1.78	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:76:PRO:HG2	38:DD:98:VAL:HG21	1.54	0.87
41:DG:5:VAL:HG12	41:DG:6:ALA:N	1.89	0.87
43:DI:92:VAL:HG11	43:DI:120:ILE:HD12	1.57	0.87
1:AA:1437:C:H42	1:AA:1464:G:H1	1.19	0.87
35:BA:1019:U:H3	35:BA:1142(A):A:H62	1.22	0.87
35:BA:1902:C:H1'	38:BD:244:ARG:HD3	1.55	0.87
35:BA:2733:A:N1	39:BE:203:LYS:HA	1.89	0.87
4:CD:100:ARG:HH12	4:CD:137:SER:HA	1.38	0.87
35:DA:1255:U:H5'	35:DA:1256:G:H5''	1.55	0.87
35:DA:528:A:H2	35:DA:2043:C:H5'	1.38	0.87
35:DA:2733:A:N1	39:DE:203:LYS:HA	1.89	0.87
1:AA:382:A:H2'	1:AA:383:A:C8	2.10	0.87
16:AP:49:LEU:HD12	16:AP:50:LYS:N	1.90	0.87
35:BA:491:G:H2'	35:BA:492:A:H8	1.38	0.87
29:D3:56:VAL:HG12	29:D3:57:GLU:H	1.38	0.87
35:DA:1114:G:H2'	35:DA:1115:G:H5''	1.57	0.87
35:DA:2206:G:N2	35:DA:2207:G:H5'	1.90	0.87
35:DA:27:G:N2	35:DA:512:G:H2'	1.90	0.87
35:DA:588:U:H2'	35:DA:589:C:C6	2.10	0.87
38:DD:53:PHE:O	38:DD:54:ARG:HG2	1.74	0.87
40:DF:114:VAL:HG23	40:DF:115:ALA:H	1.39	0.87
27:B1:25:LYS:HB2	27:B1:37:ILE:HD11	1.56	0.86
27:B1:62:VAL:HG21	27:B1:67:ILE:HA	1.57	0.86
1:CA:939:G:H5''	7:CG:102:ARG:NH2	1.90	0.86
13:CM:118:ALA:HB1	13:CM:119:GLY:N	1.89	0.86
16:CP:49:LEU:HD12	16:CP:50:LYS:N	1.90	0.86
34:D8:30:ARG:HH21	46:DP:62:LEU:HB2	1.40	0.86
41:DG:95:ARG:HA	41:DG:95:ARG:HH11	1.40	0.86
54:DX:25:LYS:HZ3	54:DX:87:GLN:N	1.73	0.86
9:AI:19:LEU:O	9:AI:20:ARG:HG3	1.74	0.86
41:BG:60:LEU:O	41:BG:63:ILE:HG13	1.75	0.86
50:BT:30:VAL:HG12	50:BT:44:ASP:HA	1.57	0.86
52:BV:72:VAL:HA	52:BV:88:ARG:HH12	1.38	0.86
55:BY:8:LYS:HD2	55:BY:8:LYS:H	1.39	0.86
1:CA:1116:C:H2'	1:CA:1117:G:H5''	1.54	0.86
41:DG:111:LEU:HA	41:DG:114:ILE:HD11	1.57	0.86
8:AH:12:ARG:HH12	8:AH:27:PRO:HD3	1.40	0.86
8:AH:83:ILE:HD13	8:AH:137:VAL:HG22	1.55	0.86
27:B1:13:ILE:HG13	27:B1:14:VAL:H	1.40	0.86
41:BG:53:LEU:HD22	41:BG:53:LEU:H	1.40	0.86
46:BP:47:ASP:HB3	46:BP:48:PRO:C	1.93	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:BX:72:LYS:HG3	54:BX:74:PRO:CD	2.02	0.86
1:CA:266:G:H5''	1:CA:268:C:H41	1.39	0.86
4:CD:100:ARG:NH1	4:CD:137:SER:HA	1.89	0.86
12:CL:47:LYS:HB3	12:CL:48:PRO:HD3	1.57	0.86
35:DA:2704:C:H2'	35:DA:2705:A:H8	1.40	0.86
54:DX:55:ASN:C	54:DX:77:LYS:HG3	1.95	0.86
7:AG:60:LYS:HA	7:AG:63:LYS:HB3	1.57	0.86
12:AL:6:THR:CG2	12:AL:9:GLN:HG3	2.05	0.86
23:AW:19:G:H3'	23:AW:20:G:H5''	1.54	0.86
35:BA:2206:G:N2	35:BA:2207:G:H5'	1.91	0.86
50:BT:50:ILE:HG23	50:BT:99:LEU:CD1	2.06	0.86
3:CC:150:LYS:HB3	3:CC:201:TYR:HB2	1.57	0.86
6:CF:7:ASN:O	6:CF:8:ILE:HG13	1.75	0.86
27:D1:77:ALA:C	27:D1:78:LYS:HD2	1.94	0.86
38:DD:265:PRO:HG2	38:DD:266:SER:H	1.39	0.86
46:DP:127:ALA:HB3	46:DP:130:PHE:CE2	2.11	0.86
1:AA:736:C:H2'	1:AA:737:A:C8	2.10	0.86
41:BG:4:ASP:HA	41:BG:8:LYS:HD2	1.57	0.86
43:BI:91:SER:CB	43:BI:119:PRO:HB2	2.04	0.86
44:BN:9:VAL:HG12	44:BN:10:GLU:H	1.41	0.86
51:BU:64:ARG:CB	51:BU:64:ARG:HH21	1.87	0.86
1:CA:34:C:H2'	1:CA:35:G:H8	1.40	0.86
2:CB:67:THR:HG21	2:CB:155:LEU:HG	1.57	0.86
6:CF:68:PRO:HG3	6:CF:71:ARG:HH21	1.40	0.86
35:DA:678:C:H2'	35:DA:679:C:H6	1.39	0.86
38:DD:8:PRO:CB	38:DD:14:ARG:HD3	2.06	0.86
44:DN:74:ARG:NH2	44:DN:101:HIS:HB3	1.90	0.86
51:DU:3:ARG:HH11	51:DU:3:ARG:HG2	1.38	0.86
1:AA:1169:A:H2'	1:AA:1170:A:H8	1.38	0.86
2:AB:97:TRP:CH2	2:AB:173:ALA:HA	2.11	0.86
7:AG:69:VAL:HA	7:AG:138:LYS:HD2	1.56	0.86
8:AH:11:THR:HG23	8:AH:14:ARG:HH12	1.39	0.86
18:AR:53:ARG:HH22	18:AR:60:ALA:N	1.74	0.86
28:B2:37:PHE:CE2	28:B2:40:SER:HA	2.11	0.86
35:BA:708:C:H42	35:BA:723:G:H1	1.23	0.86
1:CA:920:U:H1'	1:CA:1080:A:C2	2.11	0.86
3:CC:15:THR:HG22	3:CC:16:ARG:NH1	1.90	0.86
12:CL:41:ARG:HH11	12:CL:41:ARG:HB3	1.40	0.86
18:CR:53:ARG:HH22	18:CR:60:ALA:N	1.72	0.86
35:DA:1986:A:H3'	35:DA:1987:G:H5''	1.56	0.86
46:DP:77:ARG:HB2	46:DP:78:PRO:HD2	1.55	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:31:LEU:HB3	55:DY:32:PRO:CA	2.04	0.86
1:AA:1325:C:H2'	1:AA:1325:C:O2	1.76	0.86
6:AF:45:LEU:HD12	6:AF:46:ARG:H	1.39	0.86
1:AA:643:C:H5'	8:AH:31:PHE:CD1	2.09	0.86
13:AM:19:LEU:HA	13:AM:22:ILE:HD13	1.57	0.86
27:B1:22:GLY:HA2	27:B1:39:LYS:HB3	1.58	0.86
38:BD:53:PHE:O	38:BD:54:ARG:HG2	1.76	0.86
47:BQ:75:THR:CA	47:BQ:88:GLY:HA3	2.06	0.86
51:BU:92:ARG:HD3	51:BU:94:ASN:HD22	1.41	0.86
1:CA:392:G:H2'	1:CA:393:A:H8	1.40	0.86
1:CA:434:U:H2'	1:CA:435:C:C6	2.10	0.86
1:CA:489:C:H2'	1:CA:490:G:H8	1.38	0.86
2:CB:97:TRP:CH2	2:CB:173:ALA:HA	2.09	0.86
2:CB:42:ILE:HD11	2:CB:202:PRO:HB2	1.58	0.86
34:D8:59:LYS:C	34:D8:61:LEU:H	1.76	0.86
35:DA:176:G:O2'	35:DA:177:G:H5'	1.75	0.86
35:DA:1884:A:C2'	35:DA:1885:A:H5''	2.05	0.86
35:DA:1925:C:O2'	35:DA:1926:U:H5'	1.75	0.86
38:DD:27:THR:HG23	38:DD:28:GLU:N	1.91	0.86
39:DE:194:GLY:O	39:DE:196:VAL:HG23	1.75	0.86
47:DQ:52:VAL:HG13	47:DQ:53:ALA:N	1.90	0.86
46:BP:7:ARG:HB3	46:BP:8:PRO:HD3	1.56	0.86
8:CH:20:TYR:HA	8:CH:65:TYR:CE2	2.10	0.86
31:D5:47:PRO:HA	31:D5:57:VAL:HG11	1.57	0.86
38:DD:226:MET:CE	38:DD:230:ASP:HB3	2.06	0.86
43:DI:102:SER:HB2	43:DI:109:ILE:HG12	1.58	0.86
52:DV:72:VAL:HA	52:DV:88:ARG:HH12	1.39	0.86
42:BH:128:PRO:HG2	42:BH:129:THR:HG23	1.58	0.86
48:BR:9:LYS:O	48:BR:10:LEU:HG	1.76	0.86
8:CH:12:ARG:HH12	8:CH:27:PRO:HD3	1.40	0.86
34:D8:52:LYS:N	34:D8:53:PRO:HD2	1.90	0.86
35:DA:271(U):G:H2'	35:DA:271(V):G:H8	1.39	0.86
47:DQ:20:ALA:HB2	47:DQ:99:PRO:HG2	1.56	0.86
54:DX:25:LYS:HZ3	54:DX:87:GLN:H	0.90	0.86
1:AA:1513:A:H2'	1:AA:1514:C:C6	2.10	0.86
1:AA:673:G:H2'	1:AA:674:G:C8	2.11	0.86
8:AH:103:VAL:HG21	8:AH:109:ILE:O	1.75	0.86
8:AH:82:HIS:HD2	8:AH:138:TRP:HE1	1.19	0.86
41:BG:170:ARG:HH22	41:BG:182:LYS:HE2	1.39	0.86
44:BN:89:LYS:O	44:BN:93:THR:HG22	1.76	0.86
45:BO:111:PHE:HB3	45:BO:114:ILE:HD13	1.55	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:23:ARG:O	50:BT:25:GLY:N	2.08	0.86
54:BX:65:ARG:CZ	54:BX:66:LEU:H	1.89	0.86
56:BZ:102:LEU:HD11	56:BZ:124:ILE:HG23	1.58	0.86
8:CH:103:VAL:HG21	8:CH:109:ILE:O	1.76	0.86
8:CH:6:ILE:HG21	8:CH:85:ARG:NH1	1.89	0.86
13:CM:19:LEU:HA	13:CM:22:ILE:HD13	1.58	0.86
35:DA:836:G:H2'	35:DA:837:C:C6	2.11	0.86
41:DG:15:VAL:HG13	41:DG:175:LEU:HB3	1.55	0.86
54:DX:62:LYS:HB2	54:DX:68:ARG:HB2	1.58	0.86
27:B1:94:LEU:HD22	27:B1:95:LEU:N	1.90	0.85
35:BA:740:U:H2'	35:BA:741:G:C8	2.11	0.85
48:BR:24:GLN:HB2	48:BR:44:LEU:HD21	1.57	0.85
14:CN:8:GLU:HB2	14:CN:12:ARG:HH11	1.38	0.85
16:CP:20:VAL:HG21	16:CP:32:TYR:CB	2.05	0.85
25:CY:129:ILE:HA	25:CY:132:ILE:HD11	1.55	0.85
25:CY:3:LEU:N	25:CY:3:LEU:HD12	1.90	0.85
35:DA:2831:G:H1'	35:DA:2883:A:H2'	1.57	0.85
52:DV:18:LEU:HD22	52:DV:19:LYS:N	1.90	0.85
9:AI:114:TYR:H	9:AI:114:TYR:HD2	1.21	0.85
47:BQ:8:LYS:HG3	47:BQ:9:TYR:H	1.39	0.85
35:BA:106:C:H1'	55:BY:2:ARG:HH21	1.37	0.85
55:BY:45:VAL:HG13	55:BY:62:GLU:OE2	1.75	0.85
1:CA:194:C:C2'	1:CA:195:A:H5''	2.06	0.85
2:CB:75:LYS:HA	2:CB:78:GLN:NE2	1.91	0.85
19:CS:70:LYS:HB3	19:CS:70:LYS:HZ3	1.40	0.85
27:D1:60:PHE:HD1	27:D1:70:VAL:HG13	1.40	0.85
39:DE:163:GLU:O	39:DE:165:VAL:HG23	1.75	0.85
46:DP:45:LEU:HD22	46:DP:48:PRO:HB3	1.58	0.85
52:DV:32:THR:HG22	52:DV:33:VAL:H	1.41	0.85
55:DY:8:LYS:H	55:DY:8:LYS:HD2	1.41	0.85
56:DZ:163:LEU:HD23	56:DZ:163:LEU:H	1.40	0.85
12:AL:47:LYS:HB3	12:AL:48:PRO:HD3	1.55	0.85
35:BA:17:G:H4'	51:BU:25:TRP:CH2	2.11	0.85
35:BA:528:A:H2	35:BA:2043:C:H5'	1.41	0.85
35:BA:689:A:H2'	35:BA:690:G:C8	2.11	0.85
47:BQ:127:ILE:HG22	47:BQ:128:LYS:H	1.40	0.85
9:CI:19:LEU:O	9:CI:20:ARG:HG3	1.76	0.85
36:DB:7:G:H4'	49:DS:29:PHE:CD2	2.11	0.85
38:DD:77:ALA:HB2	38:DD:97:TYR:HA	1.58	0.85
47:DQ:81:VAL:HG23	47:DQ:82:ARG:NH1	1.90	0.85
1:AA:1458:G:H2'	1:AA:1459:C:C6	2.11	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AE:8:GLU:HA	5:AE:34:VAL:HG22	1.57	0.85
18:AR:85:LEU:HG	18:AR:86:VAL:H	1.41	0.85
27:B1:28:GLY:C	27:B1:30:VAL:H	1.79	0.85
35:BA:1884:A:C2'	35:BA:1885:A:H5''	2.06	0.85
40:BF:110:LEU:HA	40:BF:183:VAL:HG12	1.58	0.85
2:CB:171:ALA:HA	2:CB:174:VAL:CG2	2.05	0.85
35:DA:377:C:H2'	35:DA:378:C:C6	2.11	0.85
35:DA:676:A:H2	35:DA:802:A:H61	1.22	0.85
38:DD:35:LYS:HE3	38:DD:64:ILE:C	1.97	0.85
1:AA:543:C:H2'	1:AA:544:G:H8	1.41	0.85
1:AA:939:G:H5''	7:AG:102:ARG:NH2	1.90	0.85
10:AJ:50:ILE:HD11	14:AN:41:ARG:HD2	1.58	0.85
38:BD:131:LEU:HA	38:BD:190:TYR:CE2	2.11	0.85
1:CA:1424:C:H2'	1:CA:1425:U:H6	1.41	0.85
1:CA:678:U:H2'	1:CA:679:C:C6	2.12	0.85
39:DE:30:PRO:HD3	39:DE:180:ASN:HD21	1.42	0.85
47:DQ:39:PRO:CB	47:DQ:99:PRO:HD3	2.03	0.85
45:DO:76:ALA:HB3	50:DT:75:ILE:HB	1.59	0.85
6:AF:7:ASN:O	6:AF:8:ILE:HG13	1.75	0.85
35:BA:1590:U:C2'	35:BA:1591:G:H5''	2.06	0.85
17:CQ:40:LYS:HD2	17:CQ:42:TYR:CZ	2.11	0.85
35:DA:2262:U:C2'	35:DA:2263:C:H5''	2.06	0.85
35:DA:2262:U:H2'	35:DA:2263:C:H5''	1.59	0.85
1:AA:1410:G:H2'	1:AA:1411:C:C6	2.11	0.85
1:AA:1507:A:H2'	1:AA:1508:G:C8	2.12	0.85
2:AB:171:ALA:HA	2:AB:174:VAL:HG23	1.56	0.85
38:BD:161:THR:HG1	38:BD:196:VAL:HG21	1.40	0.85
54:BX:56:THR:N	54:BX:77:LYS:HG3	1.90	0.85
56:BZ:166:SER:HB2	56:BZ:167:PRO:HA	1.55	0.85
23:CW:19:G:H3'	23:CW:20:G:H5''	1.58	0.85
34:D8:14:VAL:HG11	34:D8:22:VAL:HG13	1.56	0.85
52:DV:71:LEU:HD13	52:DV:72:VAL:H	1.39	0.85
1:AA:17:U:H1'	1:AA:1079:G:H21	1.40	0.85
1:AA:1391:U:H2'	1:AA:1392:G:C8	2.12	0.85
35:BA:1582:C:HO2'	35:BA:1586:A:H8	1.19	0.85
38:BD:35:LYS:HE3	38:BD:64:ILE:C	1.96	0.85
47:BQ:87:LYS:O	47:BQ:87:LYS:HG3	1.77	0.85
47:BQ:20:ALA:HB2	47:BQ:99:PRO:HG2	1.57	0.85
52:BV:29:PRO:HD2	52:BV:32:THR:HG1	1.42	0.85
2:CB:222:ILE:HG23	2:CB:223:ILE:HG13	1.59	0.85
35:DA:1779:U:C5	35:DA:1784:A:N7	2.44	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:97:C:H2'	35:DA:98:G:H8	1.41	0.85
1:AA:194:C:C2'	1:AA:195:A:H5''	2.07	0.85
1:AA:932:C:H5''	7:AG:3:ARG:HD2	1.55	0.85
3:AC:148:GLY:HA3	3:AC:203:PHE:HB3	1.59	0.85
27:B1:19:GLN:HE21	35:BA:379:G:H21	1.22	0.85
38:BD:186:HIS:HD2	38:BD:188:GLU:H	1.25	0.85
38:BD:27:THR:HG23	38:BD:28:GLU:N	1.89	0.85
51:BU:34:LYS:HA	51:BU:34:LYS:HE2	1.59	0.85
1:CA:198:G:H2'	1:CA:199:G:H8	1.41	0.85
35:DA:729:G:O2'	35:DA:763:G:H4'	1.77	0.85
47:DQ:68:ILE:HD13	47:DQ:68:ILE:H	1.41	0.85
55:DY:14:LEU:HD12	55:DY:15:VAL:H	1.42	0.85
40:BF:155:LEU:HB2	40:BF:189:THR:HG21	1.59	0.85
51:BU:66:ASN:HD21	51:BU:70:ARG:HH21	0.85	0.85
54:BX:77:LYS:HE3	54:BX:78:LYS:H	1.41	0.85
55:BY:97:ARG:O	55:BY:97:ARG:HG3	1.77	0.85
5:CE:39:GLY:CA	5:CE:69:VAL:HB	2.07	0.85
29:D3:8:LEU:HA	29:D3:54:VAL:HG22	1.59	0.85
35:DA:2716:U:H2'	35:DA:2717:G:H8	1.42	0.85
18:AR:43:PHE:HA	18:AR:51:LEU:HD12	1.57	0.84
35:BA:1577:C:H2'	35:BA:1578:U:C6	2.12	0.84
35:BA:377:C:H2'	35:BA:378:C:C6	2.11	0.84
35:BA:942:G:H5'	46:BP:35:HIS:HB2	1.58	0.84
39:BE:52:LEU:HB2	39:BE:76:ARG:HB2	1.57	0.84
35:BA:2276:G:H5'	47:BQ:84:GLY:HA2	1.58	0.84
41:DG:135:LEU:HD21	41:DG:140:ILE:HD11	1.57	0.84
45:DO:31:LYS:C	45:DO:32:TYR:HD1	1.81	0.84
45:DO:69:ILE:HD12	45:DO:69:ILE:H	1.41	0.84
4:AD:155:LEU:O	4:AD:159:ARG:HG2	1.77	0.84
12:AL:41:ARG:HB3	12:AL:41:ARG:HH11	1.42	0.84
7:AG:148:ASN:HD21	23:AW:41:C:H4'	1.41	0.84
35:BA:676:A:H2	35:BA:802:A:H61	1.24	0.84
48:BR:60:LEU:HD23	48:BR:61:HIS:H	1.41	0.84
1:CA:1492:A:H2'	1:CA:1493:A:H8	1.41	0.84
2:CB:165:VAL:HG23	2:CB:166:ASP:N	1.87	0.84
17:CQ:74:LEU:HD12	17:CQ:75:ARG:HG2	1.59	0.84
27:D1:58:ILE:CD1	27:D1:59:THR:H	1.89	0.84
40:DF:129:PHE:HA	40:DF:142:TRP:HE1	1.41	0.84
45:DO:79:PHE:HE2	45:DO:101:PRO:HB2	1.41	0.84
3:AC:156:ARG:NH2	3:AC:161:GLU:HA	1.91	0.84
31:B5:47:PRO:HA	31:B5:57:VAL:HG11	1.58	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:B6:20:ASN:ND2	32:B6:21:TYR:H	1.74	0.84
35:BA:272(D):G:H1	35:BA:364:C:H42	1.25	0.84
45:BO:79:PHE:HE2	45:BO:101:PRO:HB2	1.42	0.84
45:BO:76:ALA:HB3	50:BT:75:ILE:HB	1.59	0.84
25:CY:68:VAL:O	25:CY:98:ALA:HA	1.77	0.84
39:DE:51:PHE:N	39:DE:74:PRO:HG3	1.91	0.84
56:DZ:110:GLY:C	56:DZ:112:ARG:H	1.76	0.84
8:AH:129:VAL:HG23	8:AH:130:GLY:H	1.39	0.84
1:AA:237:C:H4'	17:AQ:25:ARG:HH12	1.41	0.84
28:B2:53:LEU:HA	28:B2:56:GLN:HG2	1.59	0.84
51:BU:17:ILE:HA	51:BU:20:LEU:HD23	1.59	0.84
1:CA:1325:C:H2'	1:CA:1325:C:O2	1.75	0.84
1:CA:1508:G:H2'	1:CA:1509:C:H6	1.40	0.84
16:CP:49:LEU:HD11	16:CP:51:VAL:HG23	1.59	0.84
27:D1:85:LEU:HB2	27:D1:87:PRO:HD3	1.57	0.84
34:D8:32:LEU:C	34:D8:34:TRP:N	2.30	0.84
35:DA:1301:A:O2'	35:DA:1302:A:H2'	1.77	0.84
42:DH:26:VAL:O	42:DH:32:GLU:HA	1.77	0.84
47:DQ:127:ILE:HG22	47:DQ:128:LYS:H	1.42	0.84
48:DR:9:LYS:O	48:DR:10:LEU:HG	1.77	0.84
48:DR:48:VAL:O	48:DR:51:LEU:HB2	1.76	0.84
34:B8:32:LEU:C	34:B8:34:TRP:N	2.27	0.84
1:CA:1057:G:H5''	3:CC:154:SER:HB2	1.58	0.84
8:CH:83:ILE:HD13	8:CH:137:VAL:HG22	1.57	0.84
25:CY:41:LEU:H	25:CY:41:LEU:HD12	1.41	0.84
28:D2:27:GLU:O	28:D2:29:LYS:N	2.10	0.84
31:D5:2:ALA:HA	35:DA:2015:A:H1'	1.58	0.84
35:DA:2115:G:H4'	35:DA:2166:G:H22	1.42	0.84
41:DG:29:TRP:C	41:DG:31:VAL:H	1.77	0.84
49:DS:61:ASN:HD22	49:DS:62:LYS:HE3	1.42	0.84
1:AA:434:U:H2'	1:AA:435:C:C6	2.11	0.84
1:AA:551:U:H2'	1:AA:552:U:H6	1.43	0.84
11:AK:22:HIS:O	11:AK:28:THR:HG23	1.78	0.84
11:AK:29:ILE:HG22	11:AK:44:SER:HB3	1.59	0.84
27:B1:73:LEU:HA	27:B1:76:ARG:NH1	1.92	0.84
35:BA:2115:G:H4'	35:BA:2166:G:H22	1.43	0.84
35:BA:2704:C:H2'	35:BA:2705:A:H8	1.43	0.84
35:BA:678:C:H2'	35:BA:679:C:H6	1.41	0.84
37:BC:59:ARG:HB2	37:BC:62:VAL:HG22	1.58	0.84
38:BD:226:MET:CE	38:BD:230:ASP:HB3	2.07	0.84
42:BH:158:HIS:CD2	42:BH:170:ARG:HA	2.13	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BN:16:ILE:HG23	44:BN:54:VAL:HG22	1.58	0.84
44:BN:41:ASP:C	51:BU:64:ARG:NH1	2.31	0.84
35:BA:389:G:N2	46:BP:71:VAL:HG11	1.92	0.84
1:CA:1057:G:H5''	3:CC:154:SER:CB	2.07	0.84
3:CC:156:ARG:NH2	3:CC:161:GLU:HA	1.92	0.84
25:CY:18:LEU:HG	25:CY:19:GLU:H	1.42	0.84
26:D0:25:ARG:HA	26:D0:29:GLN:HE22	1.41	0.84
35:DA:1582:C:HO2'	35:DA:1586:A:H8	1.22	0.84
36:DB:20:C:H2'	36:DB:21:G:H5''	1.60	0.84
38:DD:144:ALA:HB3	38:DD:192:THR:HG22	1.57	0.84
38:DD:76:PRO:HA	38:DD:118:VAL:HB	1.58	0.84
40:DF:53:THR:O	40:DF:57:VAL:HG23	1.78	0.84
45:DO:111:PHE:HB3	45:DO:114:ILE:HD13	1.57	0.84
55:DY:28:LYS:HZ1	55:DY:30:VAL:HA	1.40	0.84
2:AB:47:THR:HG23	2:AB:202:PRO:HG2	1.59	0.84
5:AE:90:VAL:HG23	5:AE:121:LYS:HB3	1.60	0.84
35:BA:1484:G:H2'	35:BA:1485:G:H5''	1.59	0.84
35:BA:208:C:H2'	35:BA:209:C:H6	1.41	0.84
39:BE:194:GLY:O	39:BE:196:VAL:HG23	1.76	0.84
41:BG:7:LEU:HA	41:BG:10:LYS:HB2	1.59	0.84
42:BH:136:ILE:HD12	42:BH:136:ILE:H	1.42	0.84
47:BQ:68:ILE:H	47:BQ:68:ILE:HD13	1.40	0.84
51:BU:3:ARG:HG2	51:BU:3:ARG:HH11	1.40	0.84
55:BY:86:ARG:HB3	55:BY:88:LYS:HZ2	1.42	0.84
47:BQ:25:ASP:HA	56:BZ:78:LYS:NZ	1.93	0.84
1:CA:551:U:H2'	1:CA:552:U:H6	1.43	0.84
1:CA:683:G:H2'	1:CA:684:A:H8	1.43	0.84
2:CB:36:ARG:CZ	2:CB:37:ASN:H	1.91	0.84
27:D1:33:LYS:HG2	27:D1:34:THR:N	1.92	0.84
32:D6:20:ASN:ND2	32:D6:21:TYR:H	1.76	0.84
35:DA:2348:U:C3'	35:DA:2349:G:H5''	2.08	0.84
43:DI:111:PRO:O	43:DI:112:LYS:HG3	1.77	0.84
56:DZ:56:VAL:CA	56:DZ:70:LEU:HD21	1.99	0.84
1:AA:34:C:H2'	1:AA:35:G:H8	1.41	0.84
5:AE:39:GLY:CA	5:AE:69:VAL:HB	2.08	0.84
9:AI:95:LYS:HD3	9:AI:96:LEU:H	1.41	0.84
35:BA:1114:G:H2'	35:BA:1115:G:H5''	1.57	0.84
35:BA:1925:C:O2'	35:BA:1926:U:H5'	1.77	0.84
35:BA:2262:U:C2'	35:BA:2263:C:H5''	2.08	0.84
35:BA:271(U):G:H2'	35:BA:271(V):G:H8	1.43	0.84
35:BA:610:G:H22	35:BA:619:G:H1'	1.43	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:131:ALA:HB3	39:BE:134:ILE:HD11	1.59	0.84
47:BQ:81:VAL:HG23	47:BQ:82:ARG:NH1	1.93	0.84
35:BA:483:A:C1'	55:BY:47:LYS:HG2	2.08	0.84
28:D2:14:ARG:CG	28:D2:15:LYS:H	1.87	0.84
35:DA:2850:A:H2'	35:DA:2851:A:H8	1.42	0.84
39:DE:167:VAL:HG22	39:DE:168:MET:H	1.43	0.84
42:DH:70:THR:HG22	42:DH:74:ASN:ND2	1.92	0.84
2:AB:208:ILE:HA	2:AB:211:ILE:HD12	1.60	0.84
25:AY:29:ARG:HB3	25:AY:32:ARG:CZ	2.07	0.84
31:B5:40:LYS:CE	31:B5:46:CYS:H	1.90	0.84
35:BA:1409:C:H2'	35:BA:1410:G:C8	2.12	0.84
40:BF:53:THR:O	40:BF:57:VAL:HG23	1.78	0.84
44:BN:41:ASP:N	51:BU:64:ARG:HH12	1.75	0.84
48:BR:99:LYS:O	48:BR:100:LEU:HD22	1.78	0.84
11:CK:29:ILE:HG22	11:CK:44:SER:CB	2.07	0.84
39:DE:116:VAL:HG22	39:DE:122:PHE:HB2	1.57	0.84
41:DG:178:PHE:HB3	41:DG:180:PHE:HE1	1.43	0.84
35:DA:666:G:H4'	46:DP:49:ARG:NH2	1.93	0.84
46:DP:62:LEU:CD1	46:DP:62:LEU:H	1.91	0.84
48:DR:95:THR:HA	48:DR:117:VAL:HB	1.59	0.84
2:AB:75:LYS:HA	2:AB:78:GLN:NE2	1.92	0.84
27:B1:73:LEU:HD21	27:B1:94:LEU:HD23	1.57	0.84
35:BA:176:G:O2'	35:BA:177:G:H5'	1.78	0.84
35:BA:97:C:H2'	35:BA:98:G:H8	1.43	0.84
41:BG:172:LEU:HG	41:BG:176:LEU:HD11	1.60	0.84
1:CA:1442(A):G:H3'	1:CA:1442(B):A:C5'	2.07	0.84
1:AA:198:G:H2'	1:AA:199:G:H8	1.41	0.83
2:AB:222:ILE:HG23	2:AB:223:ILE:HG13	1.60	0.83
3:AC:24:ALA:HB3	3:AC:29:TYR:HD1	1.43	0.83
12:AL:23:LYS:O	12:AL:24:VAL:HG23	1.78	0.83
31:B5:11:THR:OG1	35:BA:1264:G:H5'	1.77	0.83
35:BA:2186:G:H2'	35:BA:2187:G:H5''	1.61	0.83
35:BA:323:G:H2'	40:BF:169:ASN:HD21	1.43	0.83
41:BG:67:LYS:H	41:BG:67:LYS:HD2	1.43	0.83
46:BP:62:LEU:CD1	46:BP:62:LEU:H	1.90	0.83
56:BZ:10:ARG:HB3	56:BZ:36:LYS:HB2	1.57	0.83
1:CA:591:U:H2'	1:CA:592:G:C8	2.12	0.83
18:CR:85:LEU:HG	18:CR:86:VAL:H	1.40	0.83
20:CT:87:LYS:O	20:CT:91:LEU:HG	1.78	0.83
27:D1:10:LYS:O	27:D1:13:ILE:HG22	1.77	0.83
35:DA:1484:G:H2'	35:DA:1485:G:H5''	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2777:G:H5''	35:DA:2778:A:C5'	2.07	0.83
38:DD:25:THR:CG2	38:DD:81:ALA:HB1	2.08	0.83
41:DG:139:LEU:HA	41:DG:144:ILE:HG12	1.60	0.83
47:DQ:75:THR:CA	47:DQ:88:GLY:HA3	2.06	0.83
1:AA:200:G:H1	1:AA:217:C:H42	1.25	0.83
4:AD:128:VAL:HG12	4:AD:129:ASN:H	1.43	0.83
40:BF:39:TRP:CD1	40:BF:101:LEU:HB2	2.12	0.83
41:BG:173:LEU:CA	41:BG:176:LEU:HD12	2.08	0.83
42:BH:156:ALA:C	42:BH:158:HIS:H	1.81	0.83
3:CC:105:GLU:HG2	3:CC:106:VAL:H	1.41	0.83
4:CD:155:LEU:O	4:CD:159:ARG:HG2	1.78	0.83
5:CE:33:VAL:HG22	5:CE:43:LEU:HD13	1.58	0.83
1:CA:878:G:H5'	8:CH:89:PRO:HG2	1.59	0.83
48:DR:2:ARG:O	48:DR:2:ARG:HD2	1.78	0.83
50:DT:30:VAL:HG12	50:DT:44:ASP:HA	1.58	0.83
35:DA:106:C:H1'	55:DY:2:ARG:HH21	1.40	0.83
1:AA:973:G:H3'	1:AA:974:A:H5''	1.60	0.83
1:AA:1057:G:H5''	3:AC:154:SER:CB	2.09	0.83
35:BA:2801(A):A:H4'	35:BA:2802:G:H2'	1.60	0.83
36:BB:20:C:H2'	36:BB:21:G:H5''	1.59	0.83
38:BD:146:GLU:HA	38:BD:153:ALA:HA	1.60	0.83
1:CA:662:G:H2'	1:CA:663:A:C8	2.12	0.83
11:CK:21:ILE:HD11	11:CK:82:VAL:HG13	1.60	0.83
35:DA:782:A:N3	38:DD:226:MET:HG2	1.93	0.83
40:DF:110:LEU:HA	40:DF:183:VAL:HG12	1.59	0.83
42:DH:70:THR:HG22	42:DH:74:ASN:HD21	1.43	0.83
46:DP:23:PRO:HB2	46:DP:33:ARG:CD	2.07	0.83
2:AB:101:MET:O	2:AB:105:PHE:HA	1.78	0.83
1:AA:1057:G:H5''	3:AC:154:SER:HB2	1.59	0.83
19:AS:48:THR:HG22	19:AS:61:TYR:HA	1.58	0.83
25:AY:76:LEU:HD11	25:AY:99:LEU:HG	1.60	0.83
35:BA:836:G:H2'	35:BA:837:C:C6	2.13	0.83
38:BD:31:LYS:HA	38:BD:31:LYS:NZ	1.94	0.83
39:BE:131:ALA:CB	39:BE:134:ILE:HD11	2.08	0.83
46:BP:121:LYS:O	46:BP:123:LEU:HD22	1.77	0.83
1:CA:1053:G:O6	1:CA:1199:U:H2'	1.78	0.83
7:CG:50:ILE:HB	7:CG:58:PRO:HD3	1.60	0.83
28:D2:44:LEU:HD23	35:DA:61:G:C5'	2.06	0.83
35:DA:2189:U:H3'	35:DA:2190:G:H5''	1.61	0.83
38:DD:146:GLU:HA	38:DD:153:ALA:HA	1.59	0.83
39:DE:29:GLY:HA3	39:DE:180:ASN:HD21	1.43	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:24:GLY:HA2	47:DQ:100:GLY:O	1.78	0.83
3:AC:150:LYS:HB3	3:AC:201:TYR:HB2	1.58	0.83
9:AI:19:LEU:HD23	9:AI:61:ALA:HA	1.58	0.83
23:AW:72:C:H4'	35:BA:1851:U:H4'	1.61	0.83
35:BA:2348:U:C3'	35:BA:2349:G:H5''	2.07	0.83
39:BE:137:HIS:HB3	39:BE:138:PRO:HD2	1.59	0.83
42:BH:70:THR:HG22	42:BH:74:ASN:HD21	1.41	0.83
1:CA:237:C:H4'	17:CQ:25:ARG:HH12	1.41	0.83
27:D1:87:PRO:HB2	27:D1:91:LYS:CE	2.09	0.83
35:DA:2820:A:H4'	48:DR:5:LYS:HE2	1.61	0.83
52:DV:22:VAL:HB	52:DV:94:LEU:HB3	1.60	0.83
53:DW:75:TYR:CE1	53:DW:104:THR:HB	2.13	0.83
1:AA:392:G:H2'	1:AA:393:A:H8	1.41	0.83
1:AA:943:U:H2'	1:AA:944:G:H8	1.43	0.83
3:AC:189:ALA:HB3	3:AC:196:LEU:HB2	1.59	0.83
35:BA:1567:A:H2'	38:BD:84:TYR:HE2	1.42	0.83
35:BA:925:C:C2'	35:BA:926:A:H5''	2.07	0.83
49:BS:55:ALA:O	49:BS:56:LEU:HB2	1.79	0.83
52:BV:22:VAL:HB	52:BV:94:LEU:HB3	1.60	0.83
25:CY:67:VAL:HG12	25:CY:100:TYR:HA	1.58	0.83
32:D6:39:TYR:HE1	35:DA:2347:C:H4'	1.44	0.83
35:DA:1007:C:O2'	44:DN:108:PRO:HA	1.78	0.83
35:DA:2781:A:C5'	35:DA:2782:G:H5'	2.08	0.83
35:DA:942:G:H5'	46:DP:35:HIS:HB2	1.61	0.83
37:DC:59:ARG:HB2	37:DC:62:VAL:HG22	1.59	0.83
38:DD:142:VAL:HG23	38:DD:193:VAL:HA	1.58	0.83
39:DE:33:VAL:CG1	39:DE:89:ASP:H	1.92	0.83
40:DF:46:ARG:HH11	40:DF:46:ARG:HA	1.43	0.83
41:DG:102:PHE:HE1	41:DG:106:LEU:HD13	1.42	0.83
41:DG:36:LYS:HD3	41:DG:95:ARG:NH2	1.92	0.83
43:DI:57:ARG:HH11	43:DI:57:ARG:HB3	1.43	0.83
54:DX:36:LYS:HD2	54:DX:36:LYS:O	1.79	0.83
56:DZ:150:LEU:HD23	56:DZ:171:ILE:CD1	2.08	0.83
47:DQ:132:VAL:HG11	56:DZ:81:ARG:NH1	1.94	0.83
35:BA:956:G:OP2	47:BQ:85:LYS:HD2	1.78	0.83
42:BH:16:SER:HB2	42:BH:27:LYS:HB2	1.61	0.83
45:BO:69:ILE:HD12	45:BO:69:ILE:H	1.41	0.83
2:CB:101:MET:O	2:CB:105:PHE:HA	1.79	0.83
1:CA:1342:C:H1'	9:CI:124:GLN:HE22	1.41	0.83
11:CK:44:SER:H	11:CK:47:VAL:CG2	1.92	0.83
31:D5:40:LYS:HE2	31:D5:46:CYS:H	1.42	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1230:C:H2'	35:DA:1231:G:C8	2.13	0.83
35:DA:1582:C:O2'	35:DA:1586:A:H8	1.61	0.83
35:DA:1971:A:H1'	38:DD:240:ALA:O	1.78	0.83
38:DD:131:LEU:HD13	38:DD:136:ILE:HG12	1.58	0.83
35:DA:323:G:H2'	40:DF:169:ASN:HD21	1.43	0.83
46:DP:112:LEU:HD22	46:DP:113:LYS:N	1.93	0.83
1:AA:662:G:H2'	1:AA:663:A:C8	2.14	0.83
16:AP:82:GLN:NE2	16:AP:82:GLN:H	1.76	0.83
35:BA:1582:C:O2'	35:BA:1586:A:H8	1.60	0.83
35:BA:2189:U:H3'	35:BA:2190:G:H5''	1.61	0.83
35:BA:484:C:H2'	35:BA:485:C:C6	2.13	0.83
35:BA:600:G:H1	35:BA:657:U:H3	1.27	0.83
38:BD:267:SER:C	38:BD:269:PHE:H	1.79	0.83
35:BA:2562:U:H1'	45:BO:23:ARG:NH1	1.92	0.83
48:BR:48:VAL:O	48:BR:51:LEU:HB2	1.79	0.83
1:CA:1240:U:H3	7:CG:30:ILE:HG22	1.43	0.83
12:CL:6:THR:CG2	12:CL:9:GLN:HG3	2.08	0.83
27:D1:46:LEU:H	27:D1:46:LEU:HD12	1.40	0.83
28:D2:29:LYS:HA	28:D2:32:LEU:HB3	1.59	0.83
35:DA:1930:G:N2	35:DA:1968:G:H2'	1.94	0.83
35:DA:2591:C:H2'	35:DA:2592:G:C8	2.14	0.83
37:DC:70:LYS:HB3	37:DC:72:VAL:HG23	1.60	0.83
44:DN:9:VAL:HG12	44:DN:10:GLU:H	1.42	0.83
44:DN:56:ASN:HA	44:DN:124:ALA:HA	1.59	0.83
54:DX:77:LYS:HE3	54:DX:78:LYS:H	1.43	0.83
1:AA:1505:G:H4'	1:AA:1506:U:H5''	1.60	0.83
12:AL:84:LEU:HD23	12:AL:85:ILE:H	1.44	0.83
35:BA:1755:A:C2	35:BA:2716:U:H1'	2.13	0.83
35:BA:2737:G:H2'	35:BA:2738:A:C8	2.13	0.83
44:BN:17:ASP:C	44:BN:19:GLU:H	1.82	0.83
47:BQ:28:ALA:HB2	47:BQ:67:ARG:HD2	1.60	0.83
54:BX:57:LEU:HB2	54:BX:76:ARG:HD2	1.59	0.83
54:BX:77:LYS:CE	54:BX:78:LYS:H	1.92	0.83
1:CA:1112:C:O2	3:CC:179:ARG:HG2	1.79	0.83
1:CA:392:G:H2'	1:CA:393:A:C8	2.14	0.83
1:CA:736:C:H2'	1:CA:737:A:C8	2.14	0.83
8:CH:26:VAL:HG22	8:CH:32:LYS:NZ	1.93	0.83
25:CY:73:GLN:O	25:CY:77:LYS:HG2	1.79	0.83
35:DA:1019:U:H2'	35:DA:1020:A:H8	1.44	0.83
35:DA:2577:A:H5''	35:DA:2578:G:H5'	1.61	0.83
35:DA:483:A:C1'	55:DY:47:LYS:HG2	2.09	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1053:G:O6	1:AA:1199:U:H2'	1.79	0.83
1:AA:1080:A:H5'	5:AE:14:ARG:HH21	1.41	0.83
6:AF:33:TYR:HD1	6:AF:75:LEU:HG	1.44	0.83
8:AH:83:ILE:HB	8:AH:137:VAL:HG13	1.61	0.83
27:B1:86:SER:C	27:B1:89:GLU:HG2	1.99	0.83
48:BR:78:LYS:O	48:BR:83:ILE:HG12	1.79	0.83
50:BT:13:ARG:HH12	50:BT:15:VAL:CG1	1.92	0.83
55:BY:28:LYS:NZ	55:BY:37:VAL:HA	1.94	0.83
1:CA:78:G:N2	1:CA:91:C:H42	1.77	0.83
1:CA:878:G:H5''	8:CH:89:PRO:HG2	1.61	0.83
27:D1:46:LEU:N	27:D1:46:LEU:HD12	1.94	0.83
28:D2:52:ASP:C	28:D2:54:LYS:H	1.79	0.83
33:D7:8:ASN:HD22	33:D7:9:ARG:N	1.76	0.83
35:DA:1409:C:H2'	35:DA:1410:G:C8	2.13	0.83
35:DA:389:G:N2	46:DP:71:VAL:HG11	1.93	0.83
3:AC:53:ALA:HB2	3:AC:115:LEU:HD21	1.61	0.82
6:AF:99:ALA:HB3	18:AR:29:PHE:HE2	1.42	0.82
8:AH:119:LEU:N	8:AH:119:LEU:HD23	1.94	0.82
1:AA:1342:C:H1'	9:AI:124:GLN:HE22	1.44	0.82
35:BA:389:G:H1	46:BP:71:VAL:CB	1.91	0.82
43:BI:37:VAL:HG13	43:BI:38:LEU:HD12	1.58	0.82
47:BQ:132:VAL:HG11	56:BZ:81:ARG:HH11	1.42	0.82
52:BV:38:LEU:HD22	52:BV:40:LEU:H	1.44	0.82
54:BX:25:LYS:HZ3	54:BX:87:GLN:H	1.22	0.82
10:CJ:50:ILE:HD11	14:CN:41:ARG:HD2	1.61	0.82
35:DA:1689:A:H62	35:DA:1698:A:H2	1.27	0.82
35:DA:2128:C:H2'	35:DA:2173:A:O2'	1.77	0.82
35:DA:2206:G:H21	35:DA:2207:G:H5'	1.44	0.82
42:DH:156:ALA:C	42:DH:158:HIS:H	1.81	0.82
43:DI:17:GLN:HG2	43:DI:18:VAL:H	1.44	0.82
35:DA:814:C:H5''	52:DV:86:GLY:HA3	1.61	0.82
55:DY:88:LYS:HZ3	55:DY:93:GLY:N	1.75	0.82
56:DZ:76:LEU:CA	56:DZ:84:GLU:HB2	2.08	0.82
1:AA:1349:A:P	9:AI:118:LYS:HZ2	2.02	0.82
1:AA:736:C:H2'	1:AA:737:A:H8	1.42	0.82
4:AD:163:GLU:O	4:AD:165:MET:N	2.12	0.82
11:AK:21:ILE:HD11	11:AK:82:VAL:HG13	1.59	0.82
27:B1:56:GLN:O	27:B1:57:GLU:HG2	1.77	0.82
27:B1:91:LYS:O	27:B1:94:LEU:HB3	1.78	0.82
28:B2:37:PHE:CZ	28:B2:40:SER:HA	2.14	0.82
45:BO:86:ILE:N	45:BO:86:ILE:HD12	1.92	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BQ:24:GLY:HA2	47:BQ:100:GLY:O	1.79	0.82
4:CD:11:LEU:C	4:CD:13:ARG:N	2.31	0.82
49:DS:83:LYS:HG2	49:DS:105:ALA:HB3	1.61	0.82
31:D5:25:LEU:HB2	53:DW:23:LEU:HD11	1.61	0.82
1:AA:376:G:OP1	16:AP:5:ARG:HB2	1.79	0.82
2:AB:36:ARG:CZ	2:AB:37:ASN:H	1.92	0.82
34:B8:14:VAL:HG11	34:B8:22:VAL:HG13	1.60	0.82
38:BD:144:ALA:HB3	38:BD:192:THR:HG22	1.59	0.82
41:BG:87:PRO:C	41:BG:88:ILE:HD12	2.00	0.82
41:BG:87:PRO:O	41:BG:88:ILE:HD12	1.79	0.82
46:BP:62:LEU:HD13	46:BP:62:LEU:H	1.44	0.82
46:BP:75:ILE:H	46:BP:75:ILE:HD12	1.41	0.82
49:BS:20:ARG:HG3	49:BS:25:ARG:HD2	1.61	0.82
50:BT:29:ARG:CB	50:BT:85:LYS:HA	2.09	0.82
6:CF:99:ALA:HB3	18:CR:29:PHE:HE2	1.44	0.82
1:CA:1329:A:P	13:CM:28:ALA:HB3	2.20	0.82
16:CP:82:GLN:H	16:CP:82:GLN:NE2	1.77	0.82
35:DA:1590:U:C2'	35:DA:1591:G:H5''	2.09	0.82
27:D1:19:GLN:HB3	35:DA:380:U:O2'	1.79	0.82
35:DA:663:G:H5''	46:DP:21:ARG:HE	1.44	0.82
36:DB:7:G:H4'	49:DS:29:PHE:CE2	2.14	0.82
41:DG:32:PRO:HB3	41:DG:172:LEU:HD22	1.62	0.82
44:DN:41:ASP:N	51:DU:64:ARG:HH12	1.77	0.82
45:DO:6:THR:HG22	45:DO:7:TYR:H	1.44	0.82
47:DQ:8:LYS:HG3	47:DQ:9:TYR:H	1.42	0.82
51:DU:90:VAL:HG12	51:DU:91:ASP:H	1.44	0.82
51:DU:90:VAL:HG13	52:DV:39:LEU:HB3	1.60	0.82
16:AP:20:VAL:HG21	16:AP:32:TYR:CB	2.07	0.82
25:AY:29:ARG:HH21	25:AY:32:ARG:HH22	1.27	0.82
35:BA:1786:A:C4	35:BA:1938:A:N6	2.48	0.82
35:BA:2476:A:C2'	35:BA:2477:C:H5''	2.09	0.82
35:BA:2598:A:C5'	38:BD:236:GLY:H	1.91	0.82
52:BV:34:GLU:HB3	52:BV:62:LEU:HD12	1.61	0.82
1:CA:382:A:H2'	1:CA:383:A:C8	2.14	0.82
5:CE:80:ILE:HD12	5:CE:138:ALA:HB1	1.60	0.82
27:D1:33:LYS:CG	27:D1:34:THR:H	1.91	0.82
35:DA:873:G:H2'	35:DA:874:G:H8	1.45	0.82
50:DT:29:ARG:CB	50:DT:85:LYS:HA	2.10	0.82
35:DA:17:G:H4'	51:DU:25:TRP:CH2	2.14	0.82
55:DY:22:GLY:O	55:DY:23:ARG:HG3	1.79	0.82
5:AE:80:ILE:HD12	5:AE:138:ALA:HB1	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1019:U:H2'	35:BA:1020:A:H8	1.41	0.82
35:BA:2639:A:C2'	35:BA:2640:G:H5''	2.09	0.82
4:CD:18:LYS:NZ	4:CD:33:MET:HB3	1.93	0.82
5:CE:36:ASP:OD2	5:CE:38:GLN:HB2	1.78	0.82
13:CM:66:LEU:HA	13:CM:70:LEU:HD12	1.61	0.82
18:CR:65:ILE:HD12	18:CR:66:LEU:N	1.94	0.82
35:DA:2639:A:C2'	35:DA:2640:G:H5''	2.07	0.82
44:DN:41:ASP:C	51:DU:64:ARG:NH1	2.32	0.82
45:DO:101:PRO:O	45:DO:102:VAL:HG13	1.80	0.82
12:AL:46:LYS:HG2	12:AL:47:LYS:H	1.44	0.82
38:BD:8:PRO:CB	38:BD:14:ARG:HD3	2.07	0.82
42:BH:144:VAL:O	42:BH:148:ILE:HG12	1.78	0.82
45:BO:1:MET:HG3	45:BO:32:TYR:CD2	2.14	0.82
51:BU:90:VAL:HG13	52:BV:39:LEU:HB3	1.58	0.82
1:CA:1391:U:H2'	1:CA:1392:G:C8	2.15	0.82
7:CG:50:ILE:O	7:CG:54:THR:HG23	1.80	0.82
18:CR:74:ARG:NE	18:CR:81:PHE:HA	1.94	0.82
31:D5:2:ALA:HB2	35:DA:2014:A:O2'	1.79	0.82
41:DG:57:ALA:HB2	41:DG:90:LEU:HD21	1.59	0.82
42:DH:128:PRO:HG2	42:DH:129:THR:HG23	1.58	0.82
1:AA:1258:G:H2'	1:AA:1259:C:H6	1.45	0.82
1:AA:253:U:H2'	1:AA:254:G:H8	1.45	0.82
1:AA:78:G:N2	1:AA:91:C:H42	1.75	0.82
8:AH:30:ARG:HB3	8:AH:30:ARG:NH1	1.95	0.82
28:B2:14:ARG:CZ	28:B2:15:LYS:H	1.93	0.82
31:B5:40:LYS:HE2	31:B5:46:CYS:H	1.43	0.82
35:BA:1773:A:H2'	35:BA:1774:C:O4'	1.80	0.82
38:BD:268:ARG:NH1	38:BD:268:ARG:HB2	1.95	0.82
49:BS:83:LYS:HG2	49:BS:105:ALA:HB3	1.61	0.82
50:BT:96:ARG:HG2	50:BT:96:ARG:HH11	1.44	0.82
56:BZ:166:SER:HB2	56:BZ:167:PRO:CA	2.10	0.82
27:D1:87:PRO:HB2	27:D1:91:LYS:NZ	1.93	0.82
31:D5:40:LYS:NZ	31:D5:45:VAL:HA	1.95	0.82
35:DA:1037:G:H1	35:DA:1118:C:H42	1.28	0.82
35:DA:484:C:H2'	35:DA:485:C:H6	1.41	0.82
45:DO:69:ILE:HD12	45:DO:77:ILE:O	1.80	0.82
35:DA:661:C:O3'	46:DP:18:ARG:HD2	1.79	0.82
54:DX:77:LYS:CE	54:DX:78:LYS:H	1.93	0.82
5:AE:36:ASP:OD2	5:AE:38:GLN:HB2	1.79	0.82
18:AR:74:ARG:NE	18:AR:81:PHE:HA	1.94	0.82
39:BE:116:VAL:HG22	39:BE:122:PHE:HB2	1.62	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BI:133:HIS:HB2	43:BI:134:PRO:HD3	1.62	0.82
51:BU:90:VAL:HG12	51:BU:91:ASP:H	1.45	0.82
3:CC:189:ALA:HB3	3:CC:196:LEU:HB2	1.61	0.82
6:CF:72:VAL:HG13	6:CF:73:ASN:H	1.44	0.82
12:CL:24:VAL:HG13	12:CL:98:TYR:CE2	2.15	0.82
31:D5:40:LYS:CE	31:D5:46:CYS:H	1.91	0.82
39:DE:173:VAL:HG12	39:DE:174:ASP:H	1.45	0.82
40:DF:101:LEU:HD12	40:DF:102:PRO:HD2	1.61	0.82
56:DZ:166:SER:HB2	56:DZ:167:PRO:HA	1.61	0.82
4:AD:176:LEU:HG	4:AD:177:ASP:H	1.44	0.82
6:AF:72:VAL:HG13	6:AF:73:ASN:H	1.45	0.82
17:AQ:74:LEU:HD12	17:AQ:75:ARG:HG2	1.60	0.82
28:B2:29:LYS:HA	28:B2:32:LEU:HD23	1.61	0.82
35:BA:1378:A:O2'	35:BA:1379:A:H5'	1.80	0.82
35:BA:2716:U:H2'	35:BA:2717:G:H8	1.45	0.82
41:BG:76:SER:HB3	41:BG:84:LYS:H	1.44	0.82
42:BH:26:VAL:O	42:BH:32:GLU:HA	1.80	0.82
43:BI:5:LEU:HD21	43:BI:19:VAL:HG12	1.61	0.82
44:BN:19:GLU:HG2	44:BN:56:ASN:O	1.78	0.82
47:BQ:24:GLY:N	47:BQ:101:ARG:HA	1.95	0.82
47:BQ:52:VAL:HG13	47:BQ:53:ALA:N	1.95	0.82
54:BX:36:LYS:HD2	54:BX:36:LYS:O	1.80	0.82
1:CA:1456:G:H2'	1:CA:1457:G:H5'	1.62	0.82
3:CC:90:GLU:O	3:CC:93:LYS:HB3	1.78	0.82
35:DA:598:G:H5'	46:DP:15:ARG:HD2	1.62	0.82
38:DD:186:HIS:HD2	38:DD:188:GLU:H	1.24	0.82
39:DE:167:VAL:O	39:DE:168:MET:HG3	1.79	0.82
50:DT:80:SER:HB3	50:DT:81:PRO:HD3	1.62	0.82
18:AR:65:ILE:HD12	18:AR:66:LEU:N	1.94	0.82
41:BG:2:PRO:O	41:BG:3:LEU:HB2	1.79	0.82
48:BR:41:ALA:HB1	48:BR:114:VAL:HG23	1.62	0.82
18:CR:25:THR:HG22	18:CR:42:ARG:NH1	1.95	0.82
20:CT:36:LEU:H	20:CT:36:LEU:HD22	1.45	0.82
41:DG:38:VAL:H	41:DG:158:ALA:HB3	1.45	0.82
42:DH:136:ILE:H	42:DH:136:ILE:HD12	1.44	0.82
46:DP:62:LEU:H	46:DP:62:LEU:HD13	1.45	0.82
1:AA:683:G:H2'	1:AA:684:A:H8	1.43	0.81
1:AA:920:U:H1'	1:AA:1080:A:C2	2.13	0.81
3:AC:52:LEU:HD23	3:AC:52:LEU:H	1.44	0.81
3:AC:90:GLU:O	3:AC:93:LYS:HB3	1.80	0.81
19:AS:36:ARG:NH2	19:AS:75:ALA:HB3	1.95	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AT:26:ASN:HB2	20:AT:71:THR:HG23	1.62	0.81
32:B6:33:LYS:HA	32:B6:33:LYS:HE2	1.62	0.81
35:BA:2850:A:H2'	35:BA:2851:A:C8	2.15	0.81
35:BA:524:U:H2'	35:BA:524:U:O2	1.80	0.81
35:BA:94(A):G:C2'	35:BA:95:G:H5''	2.10	0.81
46:BP:106:LEU:O	46:BP:107:LYS:HG2	1.78	0.81
56:BZ:127:LYS:N	56:BZ:164:ALA:HB3	1.94	0.81
25:CY:3:LEU:HD12	25:CY:3:LEU:H	1.45	0.81
40:DF:103:LYS:HG2	40:DF:106:ARG:NH2	1.95	0.81
41:DG:85:GLY:C	41:DG:87:PRO:HD3	2.00	0.81
43:DI:81:VAL:HG21	43:DI:142:VAL:HG13	1.62	0.81
48:DR:60:LEU:HD23	48:DR:61:HIS:H	1.42	0.81
47:DQ:27:VAL:CG2	56:DZ:81:ARG:HH22	1.93	0.81
32:B6:10:LEU:CD1	34:B8:36:LYS:HD3	2.08	0.81
35:BA:1570:A:H2'	35:BA:1571:A:C8	2.14	0.81
35:BA:2572:A:H2'	39:BE:144:ARG:HG3	1.62	0.81
43:BI:133:HIS:HB2	43:BI:134:PRO:CD	2.09	0.81
54:BX:65:ARG:NE	54:BX:65:ARG:HA	1.93	0.81
8:CH:11:THR:HG23	8:CH:14:ARG:HH12	1.45	0.81
35:DA:1419:A:O2'	35:DA:1420:U:H5''	1.79	0.81
35:DA:925:C:C2'	35:DA:926:A:H5''	2.09	0.81
38:DD:77:ALA:CB	38:DD:97:TYR:HA	2.11	0.81
47:DQ:24:GLY:N	47:DQ:101:ARG:HA	1.95	0.81
4:AD:5:ILE:O	4:AD:6:GLY:O	1.98	0.81
20:AT:84:LEU:O	20:AT:88:VAL:HG23	1.79	0.81
35:BA:610:G:N2	35:BA:619:G:H1'	1.96	0.81
19:CS:48:THR:HG22	19:CS:61:TYR:HA	1.59	0.81
31:D5:25:LEU:HD11	53:DW:19:LEU:HB3	1.62	0.81
42:DH:158:HIS:CD2	42:DH:170:ARG:HA	2.15	0.81
46:DP:106:LEU:O	46:DP:107:LYS:HG2	1.79	0.81
1:AA:406:G:H1	1:AA:436:C:H42	1.28	0.81
10:AJ:49:VAL:O	10:AJ:60:ARG:HB3	1.79	0.81
28:B2:30:ARG:H	28:B2:30:ARG:CD	1.90	0.81
35:BA:2777:G:H5''	35:BA:2778:A:C5'	2.09	0.81
50:BT:80:SER:HB3	50:BT:81:PRO:HD3	1.62	0.81
56:BZ:177:PRO:O	56:BZ:178:GLU:HG3	1.80	0.81
1:CA:1412:C:H2'	1:CA:1413:A:H8	1.43	0.81
12:CL:21:LYS:HD2	12:CL:21:LYS:H	1.44	0.81
12:CL:23:LYS:O	12:CL:24:VAL:HG23	1.81	0.81
12:CL:46:LYS:HG2	12:CL:47:LYS:H	1.45	0.81
12:CL:51:ALA:O	12:CL:52:LEU:HD23	1.78	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1697:G:H3'	35:DA:1698:A:H5''	1.62	0.81
35:DA:1773:A:H2'	35:DA:1774:C:O4'	1.79	0.81
41:DG:37:VAL:O	41:DG:94:LEU:HG	1.80	0.81
45:DO:22:ILE:HB	45:DO:40:VAL:HG12	1.63	0.81
46:DP:121:LYS:O	46:DP:123:LEU:HD22	1.80	0.81
27:B1:18:ILE:HG23	27:B1:42:GLN:O	1.80	0.81
34:B8:23:VAL:HG13	34:B8:47:LYS:O	1.80	0.81
35:BA:1899:G:N2	35:BA:1902:C:H41	1.78	0.81
31:B5:2:ALA:HB2	35:BA:2014:A:O2'	1.80	0.81
35:BA:2732:G:C2'	35:BA:2733:A:H5'	2.11	0.81
50:BT:13:ARG:HA	50:BT:13:ARG:CZ	2.10	0.81
1:CA:200:G:H1	1:CA:217:C:H42	1.25	0.81
3:CC:148:GLY:HA3	3:CC:203:PHE:HB3	1.61	0.81
3:CC:16:ARG:CA	3:CC:16:ARG:HH11	1.93	0.81
27:D1:64:ALA:O	27:D1:67:ILE:HG13	1.80	0.81
27:D1:88:LYS:C	27:D1:90:ILE:H	1.84	0.81
28:D2:29:LYS:HA	28:D2:32:LEU:HD23	1.62	0.81
34:D8:23:VAL:HG12	34:D8:46:ARG:HH12	1.44	0.81
35:DA:15:G:H2'	35:DA:16:G:H8	1.44	0.81
38:DD:25:THR:HB	38:DD:82:ILE:H	1.45	0.81
51:DU:88:ILE:C	51:DU:90:VAL:H	1.84	0.81
1:AA:878:G:H5''	8:AH:89:PRO:HG2	1.63	0.81
38:BD:25:THR:CG2	38:BD:81:ALA:HB1	2.11	0.81
55:BY:76:CYS:SG	55:BY:77:PRO:HD3	2.21	0.81
2:CB:96:ARG:H	2:CB:96:ARG:CD	1.89	0.81
35:DA:2220:G:H2'	35:DA:2221:G:H8	1.45	0.81
35:DA:2737:G:H2'	35:DA:2738:A:C8	2.12	0.81
35:DA:389:G:H1	46:DP:71:VAL:CB	1.94	0.81
39:DE:6:GLY:HA2	39:DE:51:PHE:CE2	2.16	0.81
47:DQ:119:ARG:HG2	47:DQ:120:ILE:HD13	1.63	0.81
49:DS:55:ALA:O	49:DS:56:LEU:HB2	1.80	0.81
52:DV:29:PRO:HD2	52:DV:32:THR:OG1	1.80	0.81
1:AA:976:G:H5'	1:AA:1358:U:O2'	1.80	0.81
35:BA:999:U:H5''	35:BA:1154:G:O6	1.80	0.81
35:BA:2128:C:H2'	35:BA:2173:A:O2'	1.80	0.81
37:BC:70:LYS:HB3	37:BC:72:VAL:HG23	1.62	0.81
43:BI:111:PRO:O	43:BI:112:LYS:HG3	1.81	0.81
54:BX:62:LYS:HB2	54:BX:68:ARG:HB2	1.61	0.81
56:BZ:56:VAL:CG2	56:BZ:70:LEU:HG	2.09	0.81
1:CA:1182:G:H4'	1:CA:1184:G:OP2	1.80	0.81
5:CE:8:GLU:HA	5:CE:34:VAL:HG22	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:119:LEU:N	8:CH:119:LEU:HD23	1.94	0.81
28:D2:14:ARG:HG2	28:D2:15:LYS:N	1.96	0.81
35:DA:2732:G:C2'	35:DA:2733:A:H5'	2.11	0.81
35:DA:460:A:H2'	35:DA:461:C:O4'	1.80	0.81
35:DA:521:G:H2'	35:DA:522:G:C8	2.16	0.81
47:DQ:28:ALA:HB2	47:DQ:67:ARG:HD2	1.62	0.81
56:DZ:48:PHE:HA	56:DZ:51:ALA:HB3	1.61	0.81
1:AA:737:A:H2'	1:AA:738:C:H6	1.45	0.81
1:AA:1112:C:O2	3:AC:179:ARG:HG2	1.81	0.81
4:AD:18:LYS:NZ	4:AD:33:MET:HB3	1.96	0.81
14:AN:16:PHE:HD2	14:AN:16:PHE:H	1.28	0.81
33:B7:9:ARG:HH12	35:BA:1309:G:H3'	1.44	0.81
35:BA:851:U:H2'	35:BA:852:G:H8	1.44	0.81
38:BD:77:ALA:HB2	38:BD:97:TYR:HA	1.61	0.81
44:BN:56:ASN:HA	44:BN:124:ALA:HA	1.62	0.81
52:BV:28:GLU:HB2	52:BV:29:PRO:CD	2.10	0.81
52:BV:83:ARG:HG2	52:BV:83:ARG:NH1	1.88	0.81
56:BZ:116:VAL:HG12	56:BZ:117:LEU:H	1.44	0.81
56:BZ:157:LEU:HD23	56:BZ:158:PRO:HD2	1.60	0.81
1:CA:253:U:H2'	1:CA:254:G:H8	1.45	0.81
3:CC:53:ALA:HB2	3:CC:115:LEU:HD21	1.62	0.81
5:CE:139:LEU:HA	5:CE:142:LEU:HD12	1.62	0.81
9:CI:79:LEU:HD21	9:CI:102:LEU:HA	1.63	0.81
9:CI:4:TYR:HB2	9:CI:19:LEU:HB2	1.62	0.81
1:CA:1372:U:H5'	9:CI:71:SER:HB3	1.62	0.81
43:DI:5:LEU:HD21	43:DI:19:VAL:HG12	1.61	0.81
56:DZ:146:ILE:HA	56:DZ:174:VAL:HB	1.62	0.81
56:DZ:73:GLN:HG2	56:DZ:74:VAL:N	1.96	0.81
1:AA:392:G:H2'	1:AA:393:A:C8	2.15	0.81
4:AD:68:TYR:HA	4:AD:114:ARG:HD3	1.63	0.81
4:AD:8:VAL:O	4:AD:10:ARG:N	2.13	0.81
12:AL:51:ALA:O	12:AL:52:LEU:HD23	1.81	0.81
31:B5:2:ALA:HA	35:BA:2015:A:H1'	1.62	0.81
33:B7:34:ARG:HB3	33:B7:42:LEU:HD23	1.63	0.81
35:BA:1846:G:H5'	35:BA:1847:A:OP2	1.81	0.81
35:BA:27:G:N2	35:BA:512:G:H2'	1.94	0.81
38:BD:76:PRO:HA	38:BD:118:VAL:HB	1.62	0.81
1:CA:41:G:H2'	1:CA:42:G:H8	1.46	0.81
2:CB:178:ARG:HH21	8:CH:74:PRO:HG3	1.46	0.81
20:CT:83:ARG:HA	20:CT:86:ARG:HB3	1.63	0.81
26:D0:40:GLN:NE2	26:D0:43:THR:HA	1.96	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:132:G:H5'	35:DA:132:G:H8	1.46	0.81
35:DA:1378:A:O2'	35:DA:1379:A:H5'	1.80	0.81
42:DH:144:VAL:O	42:DH:148:ILE:HG12	1.81	0.81
56:DZ:72:ARG:HG2	56:DZ:89:PHE:HB2	1.62	0.81
2:AB:67:THR:HG21	2:AB:155:LEU:HG	1.61	0.81
1:AA:673:G:H5''	6:AF:87:ARG:NH1	1.95	0.81
11:AK:96:ARG:HA	11:AK:99:GLN:HG2	1.61	0.81
12:AL:21:LYS:HD2	12:AL:21:LYS:H	1.46	0.81
16:AP:14:ASN:N	16:AP:15:PRO:HD3	1.95	0.81
35:BA:1133:U:O2	35:BA:1137:G:H5''	1.80	0.81
35:BA:1709:U:H2'	35:BA:1710:C:H6	1.45	0.81
40:BF:101:LEU:HD12	40:BF:102:PRO:HD2	1.63	0.81
1:CA:1191:A:H5''	3:CC:4:LYS:NZ	1.96	0.81
11:CK:43:SER:HA	11:CK:47:VAL:HG21	1.62	0.81
31:D5:11:THR:OG1	35:DA:1264:G:H5'	1.81	0.81
35:DA:1270:C:H5''	35:DA:1271:G:O5'	1.81	0.81
35:DA:1494:A:H3'	35:DA:1494:A:N3	1.94	0.81
35:DA:2128:C:H3'	35:DA:2173:A:H1'	1.62	0.81
35:DA:2712(A):A:H5'	35:DA:2713:A:OP2	1.80	0.81
37:DC:58:VAL:HG21	37:DC:166:ASP:H	1.46	0.81
38:DD:18:VAL:HG23	38:DD:211:ARG:NH2	1.95	0.81
41:DG:132:ASN:ND2	41:DG:157:ILE:HG13	1.96	0.81
25:AY:39:LEU:HA	25:AY:52:LEU:HB3	1.63	0.81
35:BA:2802:G:O2'	35:BA:2803:C:H5''	1.81	0.81
38:BD:265:PRO:HG2	38:BD:266:SER:H	1.44	0.81
39:BE:29:GLY:HA3	39:BE:180:ASN:HD21	1.43	0.81
45:BO:36:GLY:N	45:BO:62:VAL:HB	1.96	0.81
51:BU:6:THR:HG21	51:BU:10:ARG:HH21	1.46	0.81
1:CA:1191:A:H5''	3:CC:4:LYS:HZ3	1.44	0.81
4:CD:28:SER:HB3	4:CD:29:PRO:HD2	1.61	0.81
7:CG:73:MET:HG2	7:CG:90:GLU:HA	1.63	0.81
28:D2:22:GLU:HG2	54:DX:5:TYR:HB2	1.61	0.81
28:D2:56:GLN:HA	28:D2:56:GLN:NE2	1.95	0.81
35:DA:1788:C:H2'	35:DA:1789:A:H8	1.44	0.81
35:DA:2186:G:H2'	35:DA:2187:G:H5''	1.62	0.81
35:DA:2777:G:H5''	35:DA:2778:A:H5'	1.62	0.81
35:DA:585:G:H2'	35:DA:1251:C:H42	1.45	0.81
42:DH:16:SER:HB2	42:DH:27:LYS:HB2	1.62	0.81
44:DN:46:VAL:HG13	44:DN:47:ALA:N	1.95	0.81
44:DN:19:GLU:HG2	44:DN:56:ASN:O	1.81	0.81
51:DU:34:LYS:HE2	51:DU:34:LYS:HA	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:105:VAL:HG21	4:AD:126:ILE:HD13	1.63	0.80
5:AE:90:VAL:CG2	5:AE:121:LYS:HB3	2.12	0.80
13:AM:120:LYS:HE3	13:AM:120:LYS:HA	1.63	0.80
20:AT:87:LYS:O	20:AT:91:LEU:HG	1.81	0.80
25:AY:84:ARG:HE	25:AY:92:PRO:HD2	1.46	0.80
31:B5:40:LYS:NZ	31:B5:45:VAL:HA	1.97	0.80
35:BA:20:C:O2'	35:BA:21:A:H5'	1.81	0.80
35:BA:2795:G:H1	35:BA:2802:G:H1	1.28	0.80
46:BP:23:PRO:HB2	46:BP:33:ARG:HD3	1.62	0.80
55:BY:28:LYS:O	55:BY:38:ILE:HB	1.80	0.80
1:CA:1258:G:H2'	1:CA:1259:C:H6	1.44	0.80
8:CH:30:ARG:NH1	8:CH:30:ARG:HB3	1.95	0.80
8:CH:86:ILE:HG12	8:CH:135:CYS:HA	1.62	0.80
14:CN:16:PHE:H	14:CN:16:PHE:HD2	1.29	0.80
35:DA:1846:G:H5'	35:DA:1847:A:OP2	1.81	0.80
35:DA:2125:G:H21	35:DA:2173:A:N6	1.79	0.80
35:DA:2245:U:H5'	35:DA:2246:G:H5'	1.61	0.80
35:DA:2476:A:C2'	35:DA:2477:C:H5''	2.10	0.80
50:DT:22:PHE:N	50:DT:22:PHE:HD2	1.78	0.80
4:AD:108:LEU:HB3	4:AD:110:PHE:CE1	2.15	0.80
16:AP:19:ILE:HG22	16:AP:36:ILE:HG13	1.63	0.80
35:BA:1509(B):A:H2'	35:BA:1510:G:H8	1.46	0.80
35:BA:2127:G:H4'	35:BA:2128:C:OP1	1.80	0.80
41:BG:63:ILE:HD12	41:BG:64:THR:H	1.46	0.80
41:BG:73:ALA:H	41:BG:87:PRO:HD2	1.46	0.80
43:BI:81:VAL:HG21	43:BI:142:VAL:HG13	1.63	0.80
53:BW:75:TYR:CE1	53:BW:104:THR:HB	2.17	0.80
56:BZ:126:VAL:HA	56:BZ:164:ALA:HB3	1.63	0.80
1:CA:728:A:H2'	1:CA:729:A:H8	1.46	0.80
10:CJ:49:VAL:O	10:CJ:60:ARG:HB3	1.82	0.80
12:CL:84:LEU:HD23	12:CL:85:ILE:H	1.43	0.80
35:DA:1567:A:H2'	38:DD:84:TYR:HE2	1.45	0.80
35:DA:2127:G:H4'	35:DA:2128:C:OP1	1.79	0.80
35:DA:963:U:H2'	35:DA:964:C:C6	2.16	0.80
39:DE:11:MET:HB3	39:DE:24:THR:HA	1.61	0.80
39:DE:131:ALA:HB3	39:DE:134:ILE:HD11	1.63	0.80
46:DP:143:GLY:C	46:DP:145:PRO:HD3	2.01	0.80
47:DQ:137:TYR:O	47:DQ:138:ASP:HB2	1.80	0.80
47:DQ:87:LYS:HG3	47:DQ:87:LYS:O	1.80	0.80
54:DX:65:ARG:CZ	54:DX:66:LEU:H	1.95	0.80
56:DZ:48:PHE:CE2	56:DZ:71:VAL:HG21	2.15	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:49:THR:H	13:AM:52:GLU:CD	1.85	0.80
18:AR:25:THR:HG22	18:AR:42:ARG:NH1	1.96	0.80
23:AW:42:C:H2'	23:AW:43:G:C8	2.15	0.80
39:BE:163:GLU:O	39:BE:165:VAL:HG23	1.80	0.80
1:CA:674:G:H2'	1:CA:675:A:H8	1.46	0.80
3:CC:24:ALA:HB3	3:CC:29:TYR:HD1	1.46	0.80
11:CK:96:ARG:HA	11:CK:99:GLN:HG2	1.64	0.80
32:D6:33:LYS:HA	32:D6:33:LYS:HE2	1.61	0.80
35:DA:2801(A):A:H4'	35:DA:2802:G:H2'	1.61	0.80
43:DI:102:SER:HB2	43:DI:109:ILE:CG1	2.11	0.80
44:DN:16:ILE:HG23	44:DN:54:VAL:HG22	1.61	0.80
45:DO:86:ILE:HD12	45:DO:86:ILE:N	1.95	0.80
51:DU:20:LEU:H	51:DU:20:LEU:HD22	1.44	0.80
52:DV:38:LEU:HD22	52:DV:40:LEU:H	1.45	0.80
7:AG:86:GLN:HG2	23:AW:33:C:H5'	1.63	0.80
31:B5:44:THR:HG22	31:B5:45:VAL:N	1.96	0.80
35:BA:1230:C:H2'	35:BA:1231:G:H8	1.44	0.80
35:BA:1914:C:H2'	35:BA:1915:U:O4'	1.81	0.80
35:BA:2206:G:H21	35:BA:2207:G:H5'	1.44	0.80
1:CA:15:G:H2'	1:CA:16:A:H8	1.45	0.80
1:CA:973:G:H3'	1:CA:974:A:H5''	1.62	0.80
3:CC:52:LEU:HD23	3:CC:52:LEU:H	1.45	0.80
5:CE:78:HIS:HE1	5:CE:143:ARG:H	1.27	0.80
47:DQ:16:ARG:HG2	47:DQ:17:LEU:H	1.44	0.80
54:DX:56:THR:N	54:DX:77:LYS:HG3	1.95	0.80
1:AA:646:U:H2'	1:AA:647:C:C6	2.16	0.80
6:AF:75:LEU:O	6:AF:79:LEU:HG	1.81	0.80
7:AG:73:MET:HG2	7:AG:90:GLU:HA	1.62	0.80
35:BA:1697:G:H3'	35:BA:1698:A:H5''	1.61	0.80
35:BA:15:G:H2'	35:BA:16:G:H8	1.44	0.80
44:BN:46:VAL:HG13	44:BN:47:ALA:N	1.96	0.80
35:BA:661:C:O3'	46:BP:18:ARG:HD2	1.82	0.80
49:BS:25:ARG:HB3	49:BS:88:ASP:OD1	1.82	0.80
1:CA:32:A:H2'	1:CA:33:A:C8	2.16	0.80
1:CA:386:C:O2'	1:CA:387:U:H5'	1.82	0.80
13:CM:90:LEU:C	13:CM:92:HIS:H	1.81	0.80
35:DA:1791:A:H5'	38:DD:206:LEU:HD13	1.63	0.80
35:DA:1947:C:H2'	35:DA:1948:G:H8	1.46	0.80
35:DA:2802:G:O2'	35:DA:2803:C:H5''	1.82	0.80
35:DA:564:C:H2'	35:DA:565:C:C6	2.17	0.80
35:DA:94(A):G:C2'	35:DA:95:G:H5''	2.10	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:114:ILE:H	45:DO:114:ILE:CD1	1.95	0.80
53:DW:11:ARG:NH2	53:DW:98:LYS:HB3	1.96	0.80
54:DX:57:LEU:HB2	54:DX:76:ARG:HD2	1.63	0.80
1:AA:1184:G:H2'	1:AA:1185:G:H8	1.47	0.80
3:AC:16:ARG:HH11	3:AC:16:ARG:CA	1.94	0.80
4:AD:176:LEU:CG	4:AD:177:ASP:H	1.94	0.80
13:AM:23:TYR:HE1	13:AM:71:ARG:HB2	1.47	0.80
35:BA:1970:A:H5''	35:BA:1971:A:OP1	1.82	0.80
35:BA:2631:G:N2	39:BE:61:ARG:HH12	1.78	0.80
35:BA:1755:A:H2	35:BA:2716:U:H1'	1.47	0.80
39:BE:152:LYS:HB3	44:BN:78:TYR:CD1	2.15	0.80
1:CA:1489:G:H2'	1:CA:1490:C:H6	1.45	0.80
1:CA:643:C:H5'	8:CH:31:PHE:CD1	2.15	0.80
1:CA:66:G:H4'	1:CA:173:U:C5	2.16	0.80
5:CE:51:VAL:HB	5:CE:52:PRO:HD3	1.63	0.80
9:CI:95:LYS:HD3	9:CI:96:LEU:H	1.46	0.80
35:DA:1755:A:C2	35:DA:2716:U:H1'	2.17	0.80
35:DA:581:C:H2'	35:DA:582:G:H8	1.47	0.80
35:DA:742:G:H2'	35:DA:743:G:C8	2.16	0.80
39:DE:152:LYS:HB3	44:DN:78:TYR:CD1	2.17	0.80
41:DG:43:LEU:HD22	41:DG:44:GLY:N	1.95	0.80
50:DT:13:ARG:HH12	50:DT:15:VAL:CG1	1.94	0.80
52:DV:83:ARG:HG2	52:DV:83:ARG:NH1	1.92	0.80
5:AE:51:VAL:HB	5:AE:52:PRO:HD3	1.62	0.80
29:B3:8:LEU:HA	29:B3:54:VAL:HG22	1.62	0.80
35:BA:1971:A:H1'	38:BD:240:ALA:O	1.82	0.80
35:BA:521:G:H2'	35:BA:522:G:C8	2.17	0.80
35:BA:729:G:O2'	35:BA:763:G:H4'	1.82	0.80
35:BA:686:G:N2	35:BA:788:A:H61	1.80	0.80
52:BV:61:VAL:HG21	52:BV:99:ILE:HB	1.64	0.80
1:CA:1492:A:H2'	1:CA:1493:A:C8	2.15	0.80
1:CA:192:U:H4'	20:CT:103:GLY:HA2	1.64	0.80
2:CB:187:LEU:HD21	2:CB:204:ASN:O	1.79	0.80
17:CQ:19:VAL:HG23	17:CQ:44:ALA:HB3	1.62	0.80
31:D5:16:ARG:HG2	31:D5:16:ARG:HH11	1.46	0.80
35:DA:1788:C:O2'	35:DA:1789:A:H5'	1.81	0.80
48:DR:24:GLN:NE2	48:DR:36:THR:HG21	1.96	0.80
55:DY:97:ARG:O	55:DY:97:ARG:HG3	1.80	0.80
1:AA:1278:U:H5''	1:AA:1279:A:O4'	1.82	0.80
1:AA:32:A:H2'	1:AA:33:A:C8	2.16	0.80
8:AH:26:VAL:HG22	8:AH:32:LYS:NZ	1.96	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:B0:25:ARG:HA	26:B0:29:GLN:HE22	1.47	0.80
35:BA:2712(A):A:H5'	35:BA:2713:A:OP2	1.81	0.80
35:BA:30:G:H2'	35:BA:31:C:C6	2.16	0.80
38:BD:142:VAL:HG23	38:BD:193:VAL:HA	1.63	0.80
39:BE:39:PRO:HA	39:BE:43:GLY:HA2	1.63	0.80
35:BA:2636:U:H4'	39:BE:80:GLU:CD	2.02	0.80
40:BF:108:LYS:O	40:BF:111:ALA:HB3	1.80	0.80
45:BO:31:LYS:C	45:BO:32:TYR:HD1	1.85	0.80
45:BO:69:ILE:HD12	45:BO:77:ILE:O	1.79	0.80
55:BY:22:GLY:O	55:BY:23:ARG:HG3	1.82	0.80
1:CA:1278:U:H5''	1:CA:1279:A:O4'	1.82	0.80
4:CD:108:LEU:HB3	4:CD:110:PHE:CE1	2.16	0.80
8:CH:68:ARG:HG3	8:CH:69:ARG:H	1.47	0.80
32:D6:11:LEU:HG	32:D6:26:ASN:ND2	1.97	0.80
35:DA:1133:U:O2	35:DA:1137:G:H5''	1.82	0.80
35:DA:1777:U:O2'	35:DA:1778:U:H5'	1.82	0.80
45:DO:1:MET:HG3	45:DO:32:TYR:CD2	2.16	0.80
1:AA:41:G:H2'	1:AA:42:G:H8	1.46	0.80
1:AA:683:G:H2'	1:AA:684:A:C8	2.17	0.80
4:AD:28:SER:HB3	4:AD:29:PRO:HD2	1.62	0.80
12:AL:47:LYS:CB	12:AL:48:PRO:HD3	2.11	0.80
27:B1:42:GLN:HG2	27:B1:43:TYR:N	1.96	0.80
35:BA:1378:A:H4'	35:BA:1379:A:OP1	1.80	0.80
35:BA:460:A:H2'	35:BA:461:C:O4'	1.81	0.80
40:BF:53:THR:H	40:BF:56:GLU:CB	1.95	0.80
43:BI:17:GLN:HG2	43:BI:18:VAL:H	1.43	0.80
46:BP:112:LEU:HD22	46:BP:113:LYS:N	1.95	0.80
35:BA:814:C:H5''	52:BV:86:GLY:HA3	1.64	0.80
54:BX:8:ILE:H	54:BX:8:ILE:HD12	1.47	0.80
4:CD:176:LEU:HG	4:CD:177:ASP:H	1.47	0.80
18:CR:43:PHE:HA	18:CR:51:LEU:HD12	1.62	0.80
35:DA:1039:G:H1	35:DA:1116:C:N4	1.79	0.80
35:DA:1970:A:H5''	35:DA:1971:A:OP1	1.81	0.80
35:DA:2631:G:N2	39:DE:61:ARG:HH12	1.79	0.80
38:DD:267:SER:C	38:DD:269:PHE:H	1.82	0.80
40:DF:155:LEU:HB2	40:DF:189:THR:HG21	1.62	0.80
1:AA:892:A:H2'	1:AA:893:C:C6	2.17	0.80
2:AB:171:ALA:HA	2:AB:174:VAL:CG2	2.12	0.80
13:AM:23:TYR:HD1	13:AM:67:GLU:HA	1.47	0.80
19:AS:41:VAL:HB	19:AS:44:MET:HB2	1.64	0.80
26:B0:40:GLN:NE2	26:B0:43:THR:HA	1.95	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:34:ARG:HD3	33:B7:42:LEU:HA	1.63	0.80
35:BA:1947:C:H2'	35:BA:1948:G:H8	1.47	0.80
35:BA:2262:U:H2'	35:BA:2263:C:H5''	1.63	0.80
35:BA:229:A:H3'	35:BA:230:U:H5'	1.64	0.80
39:BE:11:MET:H	50:BT:8:LYS:NZ	1.80	0.80
46:BP:45:LEU:HD23	46:BP:46:LYS:H	1.46	0.80
1:CA:976:G:H5'	1:CA:1358:U:O2'	1.82	0.80
2:CB:208:ILE:HA	2:CB:211:ILE:HD12	1.62	0.80
4:CD:68:TYR:HA	4:CD:114:ARG:HD3	1.64	0.80
10:CJ:48:THR:OG1	10:CJ:62:HIS:HB3	1.82	0.80
35:DA:2393:A:H5'	46:DP:62:LEU:HB3	1.63	0.80
35:DA:292:C:H42	35:DA:348:G:H1	1.27	0.80
35:DA:30:G:H2'	35:DA:31:C:C6	2.17	0.80
36:DB:42:C:H1'	41:DG:92:VAL:HG23	1.64	0.80
38:DD:34:VAL:HG22	38:DD:35:LYS:HG3	1.63	0.80
44:DN:17:ASP:C	44:DN:19:GLU:H	1.84	0.80
45:DO:69:ILE:HD13	45:DO:77:ILE:HG23	1.64	0.80
46:DP:17:LYS:O	46:DP:17:LYS:HG2	1.80	0.80
49:DS:20:ARG:HG3	49:DS:25:ARG:HD2	1.63	0.80
1:AA:1240:U:H3	7:AG:30:ILE:HG22	1.46	0.79
1:AA:386:C:O2'	1:AA:387:U:H5'	1.82	0.79
8:AH:68:ARG:HG3	8:AH:69:ARG:H	1.47	0.79
13:AM:66:LEU:HA	13:AM:70:LEU:HD12	1.64	0.79
13:AM:90:LEU:C	13:AM:92:HIS:H	1.81	0.79
25:AY:150:SER:O	25:AY:154:THR:HG23	1.82	0.79
34:B8:22:VAL:HB	34:B8:53:PRO:CB	2.12	0.79
35:BA:1336:A:H2'	35:BA:1337:G:H8	1.47	0.79
35:BA:1456:G:H2'	35:BA:1457:A:H8	1.46	0.79
35:BA:251:A:H5''	46:BP:51:PHE:CZ	2.17	0.79
35:BA:759:G:H2'	35:BA:760:G:H8	1.45	0.79
1:CA:683:G:H2'	1:CA:684:A:C8	2.17	0.79
5:CE:19:MET:SD	5:CE:24:ARG:HG2	2.21	0.79
7:CG:79:ARG:HE	7:CG:84:ASN:ND2	1.80	0.79
35:DA:2533:A:C2'	35:DA:2534:A:H5''	2.12	0.79
35:DA:291:C:H2'	35:DA:292:C:C6	2.16	0.79
9:AI:4:TYR:HB2	9:AI:19:LEU:HB2	1.64	0.79
11:AK:44:SER:H	11:AK:47:VAL:CG2	1.95	0.79
12:AL:24:VAL:HG13	12:AL:98:TYR:CE2	2.17	0.79
1:AA:1329:A:P	13:AM:28:ALA:HB3	2.22	0.79
17:AQ:19:VAL:HG23	17:AQ:44:ALA:HB3	1.63	0.79
25:AY:38:LEU:HD12	25:AY:58:VAL:HG11	1.62	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1039:G:H1	35:BA:1116:C:N4	1.79	0.79
35:BA:2577:A:H5''	35:BA:2578:G:H5'	1.61	0.79
35:BA:2645:G:H3'	35:BA:2646:C:C5'	2.13	0.79
35:BA:291:C:H2'	35:BA:292:C:C6	2.17	0.79
35:BA:540:C:H2'	35:BA:541:C:C6	2.18	0.79
35:BA:914:C:H2'	35:BA:915:C:H5'	1.64	0.79
46:BP:84:ASN:HA	46:BP:115:LEU:O	1.83	0.79
56:BZ:138:GLU:O	56:BZ:155:LEU:HD11	1.82	0.79
4:CD:120:LEU:HD12	4:CD:120:LEU:N	1.96	0.79
8:CH:122:ARG:HA	8:CH:125:ARG:HB3	1.63	0.79
12:CL:84:LEU:HD23	12:CL:85:ILE:N	1.98	0.79
13:CM:10:PRO:CG	13:CM:18:ALA:HB1	2.12	0.79
35:DA:1456:G:H2'	35:DA:1457:A:H8	1.45	0.79
39:DE:108:SER:CB	39:DE:165:VAL:HG21	2.12	0.79
40:DF:63:LYS:NZ	40:DF:67:GLN:HB3	1.95	0.79
41:DG:141:PHE:HD1	41:DG:142:PRO:HD2	1.44	0.79
47:DQ:35:VAL:HG23	47:DQ:102:VAL:HA	1.64	0.79
1:AA:15:G:H2'	1:AA:16:A:H8	1.47	0.79
1:AA:728:A:H2'	1:AA:729:A:H8	1.47	0.79
18:AR:50:ILE:HD12	18:AR:70:ILE:HG21	1.65	0.79
20:AT:104:LEU:HD23	20:AT:105:SER:N	1.98	0.79
35:BA:1255:U:H5'	35:BA:1256:G:H5''	1.64	0.79
46:BP:30:THR:HG22	46:BP:31:ALA:N	1.96	0.79
48:BR:24:GLN:NE2	48:BR:36:THR:HG21	1.98	0.79
51:BU:20:LEU:H	51:BU:20:LEU:HD22	1.46	0.79
25:CY:68:VAL:HG23	25:CY:99:LEU:HB2	1.64	0.79
35:DA:1677:A:H2'	35:DA:1678:G:C8	2.16	0.79
35:DA:208:C:H2'	35:DA:209:C:H6	1.47	0.79
35:DA:20:C:O2'	35:DA:21:A:H5'	1.82	0.79
35:DA:8:A:C4	35:DA:9:U:H5	2.01	0.79
36:DB:74:U:C3'	36:DB:75:G:H5''	2.12	0.79
48:DR:78:LYS:O	48:DR:83:ILE:HG12	1.81	0.79
54:DX:78:LYS:HD3	54:DX:78:LYS:O	1.81	0.79
56:DZ:33:LEU:HG	56:DZ:35:ARG:H	1.47	0.79
1:AA:636:U:H2'	1:AA:637:G:C8	2.17	0.79
1:AA:735:C:O2'	1:AA:736:C:H5'	1.83	0.79
2:AB:187:LEU:HD21	2:AB:204:ASN:O	1.83	0.79
3:AC:84:ILE:HA	3:AC:87:LEU:HD12	1.64	0.79
17:AQ:40:LYS:HD2	17:AQ:42:TYR:CZ	2.18	0.79
17:AQ:69:LYS:O	17:AQ:70:ARG:HD2	1.83	0.79
31:B5:16:ARG:HH12	31:B5:17:ASP:CG	1.86	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1037:G:H1	35:BA:1118:C:H42	1.29	0.79
35:BA:2245:U:H5'	35:BA:2246:G:H5'	1.65	0.79
38:BD:25:THR:HB	38:BD:82:ILE:H	1.45	0.79
44:BN:74:ARG:CZ	44:BN:101:HIS:HB3	2.13	0.79
46:BP:101:VAL:HG13	46:BP:102:ARG:H	1.46	0.79
35:BA:626:U:O2	46:BP:105:LEU:HG	1.81	0.79
46:BP:16:ARG:HD3	46:BP:18:ARG:H	1.48	0.79
47:BQ:137:TYR:O	47:BQ:138:ASP:HB2	1.80	0.79
49:BS:27:SER:HA	49:BS:89:ARG:HD2	1.64	0.79
55:BY:14:LEU:HD12	55:BY:15:VAL:H	1.47	0.79
2:CB:22:LYS:NZ	2:CB:22:LYS:HA	1.97	0.79
4:CD:128:VAL:HG12	4:CD:129:ASN:H	1.47	0.79
8:CH:83:ILE:HB	8:CH:137:VAL:HG13	1.62	0.79
35:DA:2779:U:H1'	35:DA:2781:A:C5	2.17	0.79
35:DA:2636:U:H4'	39:DE:80:GLU:CD	2.03	0.79
40:DF:108:LYS:O	40:DF:111:ALA:HB3	1.82	0.79
43:DI:11:ASN:HD22	43:DI:12:LEU:HD22	1.46	0.79
43:DI:133:HIS:HB2	43:DI:134:PRO:CD	2.12	0.79
45:DO:2:ILE:HD11	45:DO:82:ASN:HB3	1.64	0.79
55:DY:28:LYS:NZ	55:DY:37:VAL:HA	1.98	0.79
2:AB:22:LYS:HA	2:AB:22:LYS:NZ	1.97	0.79
4:AD:11:LEU:C	4:AD:13:ARG:N	2.33	0.79
8:AH:30:ARG:CB	8:AH:30:ARG:HH11	1.94	0.79
35:BA:1255:U:C5'	35:BA:1256:G:H5''	2.12	0.79
35:BA:528:A:H2	35:BA:2043:C:C5'	1.95	0.79
35:BA:2128:C:H3'	35:BA:2173:A:H1'	1.63	0.79
35:BA:2630:G:H1'	35:BA:2894:G:H1'	1.62	0.79
35:BA:969:U:H2'	35:BA:970:C:C6	2.17	0.79
50:BT:22:PHE:HD2	50:BT:22:PHE:N	1.80	0.79
8:CH:11:THR:HA	8:CH:14:ARG:NH1	1.97	0.79
10:CJ:50:ILE:CD1	10:CJ:50:ILE:H	1.89	0.79
23:CW:17:C:H4'	23:CW:62:C:H5'	1.64	0.79
27:D1:10:LYS:HG3	27:D1:11:ARG:N	1.96	0.79
35:DA:969:U:H2'	35:DA:970:C:C6	2.18	0.79
52:DV:37:VAL:HG12	52:DV:38:LEU:H	1.46	0.79
8:AH:109:ILE:HG12	8:AH:110:ALA:N	1.98	0.79
12:AL:89:ARG:HB2	12:AL:89:ARG:NH1	1.96	0.79
32:B6:39:TYR:HE1	35:BA:2347:C:H4'	1.47	0.79
38:BD:77:ALA:CB	38:BD:97:TYR:HA	2.13	0.79
39:BE:11:MET:HB3	39:BE:24:THR:HA	1.64	0.79
43:BI:11:ASN:HD22	43:BI:12:LEU:HD22	1.46	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:45:LEU:HD22	46:BP:48:PRO:HB3	1.62	0.79
48:BR:95:THR:HA	48:BR:117:VAL:HB	1.64	0.79
49:BS:34:HIS:CD2	49:BS:53:SER:HB3	2.18	0.79
51:BU:83:LEU:HG	51:BU:88:ILE:HG21	1.63	0.79
1:CA:1489:G:H2'	1:CA:1490:C:C6	2.18	0.79
1:CA:736:C:H2'	1:CA:737:A:H8	1.48	0.79
13:CM:23:TYR:HD1	13:CM:67:GLU:HA	1.48	0.79
13:CM:49:THR:H	13:CM:52:GLU:CD	1.86	0.79
28:D2:26:ARG:NH2	54:DX:7:VAL:H	1.80	0.79
34:D8:22:VAL:HB	34:D8:53:PRO:CB	2.12	0.79
35:DA:999:U:H5''	35:DA:1154:G:O6	1.81	0.79
35:DA:2630:G:H1'	35:DA:2894:G:H1'	1.65	0.79
35:DA:635:C:H2'	35:DA:636:G:H8	1.48	0.79
40:DF:65:TRP:CH2	40:DF:75:HIS:HD2	1.99	0.79
43:DI:98:ALA:HB1	43:DI:109:ILE:CB	2.12	0.79
50:DT:13:ARG:CZ	50:DT:13:ARG:HA	2.12	0.79
52:DV:61:VAL:HG21	52:DV:99:ILE:HB	1.63	0.79
54:DX:65:ARG:HA	54:DX:65:ARG:NE	1.94	0.79
3:AC:86:VAL:O	3:AC:90:GLU:HG2	1.83	0.79
16:AP:45:THR:HG22	16:AP:47:ASP:H	1.48	0.79
35:BA:663:G:H5''	46:BP:21:ARG:HE	1.47	0.79
35:BA:782:A:N3	38:BD:226:MET:HG2	1.97	0.79
35:BA:8:A:C4	35:BA:9:U:H5	2.00	0.79
38:BD:25:THR:CB	38:BD:82:ILE:H	1.96	0.79
46:BP:143:GLY:C	46:BP:145:PRO:HD3	2.02	0.79
1:CA:1074:G:H2'	1:CA:1075:C:H6	1.48	0.79
35:DA:491:G:H2'	35:DA:492:A:C8	2.16	0.79
35:DA:600:G:H1	35:DA:657:U:H3	1.26	0.79
46:DP:101:VAL:HG13	46:DP:102:ARG:H	1.46	0.79
52:DV:61:VAL:HB	52:DV:99:ILE:H	1.46	0.79
55:DY:45:VAL:CA	55:DY:62:GLU:HG2	2.08	0.79
12:AL:70:ILE:N	12:AL:70:ILE:HD12	1.97	0.79
13:AM:10:PRO:CG	13:AM:18:ALA:HB1	2.13	0.79
31:B5:2:ALA:HB3	35:BA:747:U:C2	2.17	0.79
35:BA:1301:A:O2'	35:BA:1302:A:H2'	1.82	0.79
35:BA:1419:A:O2'	35:BA:1420:U:H5''	1.82	0.79
52:BV:29:PRO:HD2	52:BV:32:THR:OG1	1.81	0.79
1:CA:943:U:H2'	1:CA:944:G:H8	1.47	0.79
1:AA:543:C:H2'	1:AA:544:G:C8	2.18	0.79
2:AB:111:ARG:HG2	2:AB:111:ARG:HH11	1.48	0.79
5:AE:139:LEU:HA	5:AE:142:LEU:HD12	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:122:ARG:HA	8:AH:125:ARG:HB3	1.65	0.79
35:BA:1414:G:H2'	35:BA:1415:U:C6	2.18	0.79
35:BA:2000:G:HO2'	35:BA:2689:U:H5	1.28	0.79
35:BA:587:C:H3'	46:BP:33:ARG:HH12	1.45	0.79
35:BA:813:U:H2'	35:BA:814:C:C6	2.17	0.79
11:CK:121:PRO:O	11:CK:126:ARG:HB2	1.82	0.79
35:DA:2558:C:H2'	35:DA:2559:C:H6	1.48	0.79
38:DD:265:PRO:O	38:DD:267:SER:N	2.16	0.79
46:DP:16:ARG:HD3	46:DP:18:ARG:H	1.48	0.79
46:DP:59:LEU:HA	46:DP:61:ARG:CZ	2.13	0.79
49:DS:25:ARG:HB3	49:DS:88:ASP:OD1	1.83	0.79
35:DA:1754:C:H5'	50:DT:101:PHE:CD1	2.18	0.79
11:AK:29:ILE:HG22	11:AK:44:SER:CB	2.13	0.79
18:AR:56:THR:HB	18:AR:58:LEU:CD1	2.12	0.79
18:AR:87:ARG:NH1	18:AR:87:ARG:HB3	1.98	0.79
35:BA:1779:U:C5	35:BA:1784:A:N7	2.51	0.79
35:BA:2173:A:OP2	35:BA:2173:A:H3'	1.83	0.79
41:BG:131:TYR:O	41:BG:159:VAL:HG13	1.83	0.79
49:BS:61:ASN:HD22	49:BS:62:LYS:HE3	1.45	0.79
1:CA:665:A:H2'	1:CA:725:G:N2	1.98	0.79
2:CB:140:HIS:O	2:CB:143:GLU:HB2	1.83	0.79
19:CS:41:VAL:HB	19:CS:44:MET:HB2	1.65	0.79
20:CT:84:LEU:O	20:CT:88:VAL:HG23	1.83	0.79
41:DG:124:SER:HB3	41:DG:131:TYR:CE1	2.18	0.79
42:DH:67:LEU:HG	42:DH:71:LEU:HD21	1.65	0.79
35:DA:956:G:OP2	47:DQ:85:LYS:HD2	1.83	0.79
50:DT:50:ILE:HG23	50:DT:99:LEU:CD1	2.12	0.79
20:AT:36:LEU:H	20:AT:36:LEU:HD22	1.46	0.78
25:AY:73:GLN:HG3	25:AY:74:ASN:H	1.48	0.78
31:B5:45:VAL:HG22	31:B5:51:TYR:HD1	1.48	0.78
29:B3:29:ARG:NH1	35:BA:1183:G:H4'	1.98	0.78
35:BA:2125:G:H21	35:BA:2173:A:N6	1.81	0.78
40:BF:46:ARG:HA	40:BF:46:ARG:HH11	1.48	0.78
47:BQ:16:ARG:HG2	47:BQ:17:LEU:H	1.48	0.78
1:CA:1468:A:H2'	1:CA:1469:G:O4'	1.83	0.78
6:CF:75:LEU:O	6:CF:79:LEU:HG	1.83	0.78
35:DA:268:C:H2'	35:DA:268:C:O2	1.82	0.78
40:DF:84:VAL:O	40:DF:86:GLY:N	2.16	0.78
41:DG:91:ARG:C	41:DG:91:ARG:HD2	2.03	0.78
45:DO:105:GLU:N	45:DO:105:GLU:OE1	2.16	0.78
45:DO:6:THR:HG22	45:DO:7:TYR:N	1.97	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:43:THR:HA	47:DQ:94:VAL:HG12	1.65	0.78
1:AA:674:G:H2'	1:AA:675:A:H8	1.48	0.78
2:AB:55:PHE:HA	2:AB:58:ILE:HB	1.65	0.78
6:AF:10:LEU:HA	6:AF:84:ASN:O	1.82	0.78
20:AT:83:ARG:HA	20:AT:86:ARG:HB3	1.65	0.78
22:AV:30:A:H2'	22:AV:31:U:H6	1.45	0.78
25:AY:152:ASP:O	25:AY:156:ARG:HG3	1.83	0.78
34:B8:59:LYS:C	34:B8:61:LEU:H	1.83	0.78
35:BA:2777:G:H5''	35:BA:2778:A:H5'	1.63	0.78
35:BA:904:C:H2'	35:BA:904:C:O2	1.82	0.78
38:BD:172:TYR:HD1	38:BD:186:HIS:HA	1.48	0.78
41:BG:61:ALA:HA	41:BG:64:THR:CG2	2.13	0.78
46:BP:39:LYS:HD2	46:BP:40:SER:H	1.48	0.78
54:BX:12:VAL:HG11	54:BX:27:THR:HG23	1.65	0.78
1:CA:41:G:H2'	1:CA:42:G:C8	2.19	0.78
4:CD:105:VAL:HG21	4:CD:126:ILE:HD13	1.65	0.78
13:CM:23:TYR:HE1	13:CM:71:ARG:HB2	1.45	0.78
25:CY:169:ILE:O	25:CY:172:ALA:HB3	1.84	0.78
29:D3:29:ARG:NH1	35:DA:1183:G:H4'	1.98	0.78
35:DA:2795:G:H1	35:DA:2802:G:H1	1.29	0.78
35:DA:528:A:N1	35:DA:2042:A:H2'	1.99	0.78
41:DG:12:TYR:HA	41:DG:16:ARG:NH1	1.99	0.78
5:AE:150:ARG:HB2	5:AE:150:ARG:NH1	1.98	0.78
35:BA:1494:A:N3	35:BA:1494:A:H3'	1.97	0.78
35:BA:1930:G:H22	35:BA:1968:G:H2'	1.48	0.78
38:BD:231:HIS:ND1	38:BD:232:PRO:HD2	1.99	0.78
38:BD:265:PRO:O	38:BD:267:SER:N	2.16	0.78
1:CA:1458:G:H2'	1:CA:1459:C:H6	1.45	0.78
4:CD:59:ARG:HH22	4:CD:66:ARG:HH22	1.31	0.78
35:DA:2173:A:OP2	35:DA:2173:A:H3'	1.83	0.78
35:DA:2645:G:H3'	35:DA:2646:C:C5'	2.12	0.78
36:DB:42:C:O2	41:DG:93:THR:N	2.15	0.78
38:DD:264:LYS:HE2	38:DD:266:SER:O	1.82	0.78
34:D8:13:ARG:CB	46:DP:63:PRO:HA	2.14	0.78
52:DV:19:LYS:HZ2	52:DV:20:LEU:H	1.31	0.78
1:AA:1478:C:H2'	1:AA:1479:C:C6	2.18	0.78
1:AA:66:G:H4'	1:AA:173:U:C5	2.18	0.78
9:AI:79:LEU:HD21	9:AI:102:LEU:HA	1.63	0.78
1:AA:1372:U:H5''	9:AI:71:SER:HB3	1.62	0.78
35:BA:1409:C:H2'	35:BA:1410:G:H8	1.46	0.78
35:BA:1602:U:H3'	35:BA:1603:A:C5'	2.13	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1835:G:H5'	35:BA:1836:C:OP2	1.84	0.78
35:BA:2781:A:C5'	35:BA:2782:G:H5'	2.12	0.78
47:BQ:43:THR:HA	47:BQ:94:VAL:HG12	1.64	0.78
50:BT:28:VAL:HG13	50:BT:45:PHE:O	1.83	0.78
52:BV:15:GLU:HB3	52:BV:16:PRO:HD2	1.66	0.78
52:BV:61:VAL:HB	52:BV:99:ILE:H	1.48	0.78
1:CA:1346:A:H5''	9:CI:120:ARG:HH12	1.49	0.78
3:CC:84:ILE:HA	3:CC:87:LEU:HD12	1.63	0.78
5:CE:150:ARG:HB2	5:CE:150:ARG:NH1	1.98	0.78
35:DA:1414:G:H2'	35:DA:1415:U:C6	2.19	0.78
35:DA:587:C:H3'	46:DP:33:ARG:HH12	1.49	0.78
51:DU:83:LEU:HG	51:DU:88:ILE:HG21	1.63	0.78
1:AA:1299:A:H2'	1:AA:1301:U:C6	2.17	0.78
2:AB:159:PRO:C	2:AB:161:ALA:H	1.87	0.78
34:B8:50:LEU:H	34:B8:53:PRO:HG3	1.47	0.78
35:BA:2315:G:H2'	35:BA:2316:C:C6	2.18	0.78
39:BE:172:VAL:HG13	39:BE:182:LEU:HD11	1.66	0.78
43:BI:57:ARG:HH11	43:BI:57:ARG:HB3	1.48	0.78
2:CB:159:PRO:C	2:CB:161:ALA:H	1.86	0.78
13:CM:120:LYS:HE3	13:CM:120:LYS:HA	1.63	0.78
23:CW:19:G:C3'	23:CW:20:G:H5''	2.14	0.78
26:D0:27:GLU:N	26:D0:69:PHE:HE1	1.81	0.78
33:D7:9:ARG:HH12	35:DA:1309:G:H3'	1.47	0.78
35:DA:1409:C:H2'	35:DA:1410:G:H8	1.46	0.78
36:DB:7:G:H21	49:DS:38:GLN:HE22	1.28	0.78
38:DD:163:ALA:HB1	38:DD:175:LEU:HD21	1.64	0.78
35:DA:2572:A:H2'	39:DE:144:ARG:HG3	1.64	0.78
39:DE:11:MET:H	50:DT:8:LYS:NZ	1.80	0.78
51:DU:47:TYR:HD1	51:DU:50:ARG:HH22	1.32	0.78
10:AJ:48:THR:OG1	10:AJ:62:HIS:HB3	1.83	0.78
27:B1:86:SER:HA	27:B1:89:GLU:CG	2.11	0.78
35:BA:1717:G:H3'	35:BA:1718:G:H5''	1.63	0.78
35:BA:2393:A:H5'	46:BP:62:LEU:HB3	1.66	0.78
48:BR:20:LEU:HD12	48:BR:20:LEU:C	2.04	0.78
1:CA:17:U:H1'	1:CA:1079:G:H21	1.47	0.78
1:CA:972:C:H4'	10:CJ:57:LYS:CG	2.14	0.78
2:CB:55:PHE:HA	2:CB:58:ILE:HB	1.65	0.78
10:CJ:4:ILE:HG23	10:CJ:98:ILE:HG23	1.66	0.78
13:CM:15:VAL:O	13:CM:19:LEU:HD23	1.83	0.78
35:DA:1786:A:C4	35:DA:1938:A:N6	2.51	0.78
46:DP:30:THR:HG22	46:DP:31:ALA:N	1.97	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1182:G:H4'	1:AA:1184:G:OP2	1.84	0.78
8:AH:53:VAL:O	8:AH:56:LYS:HB2	1.84	0.78
13:AM:3:ARG:HA	13:AM:9:ILE:HG13	1.64	0.78
16:AP:82:GLN:HE21	16:AP:82:GLN:N	1.82	0.78
18:AR:53:ARG:HA	18:AR:56:THR:OG1	1.84	0.78
35:BA:292:C:H42	35:BA:348:G:H1	1.29	0.78
35:BA:635:C:H2'	35:BA:636:G:H8	1.48	0.78
35:BA:873:G:H2'	35:BA:874:G:H8	1.47	0.78
39:BE:116:VAL:HG21	39:BE:122:PHE:CD2	2.18	0.78
41:BG:101:ILE:HD11	41:BG:105:LYS:HE3	1.64	0.78
35:BA:1011:G:OP1	51:BU:75:ASN:HB2	1.84	0.78
52:BV:19:LYS:NZ	52:BV:20:LEU:H	1.81	0.78
52:BV:37:VAL:HG12	52:BV:38:LEU:H	1.48	0.78
55:BY:15:VAL:HG12	55:BY:16:ALA:H	1.48	0.78
1:CA:1431:C:H2'	1:CA:1432:G:O4'	1.82	0.78
5:CE:82:VAL:HG21	5:CE:138:ALA:HA	1.65	0.78
5:CE:90:VAL:C	5:CE:91:LEU:HD12	2.04	0.78
34:D8:50:LEU:H	34:D8:53:PRO:HG3	1.47	0.78
47:DQ:20:ALA:HA	47:DQ:98:LYS:HB3	1.66	0.78
54:DX:64:LYS:HG2	54:DX:65:ARG:N	1.99	0.78
56:DZ:6:LYS:HB2	56:DZ:8:TYR:CE1	2.19	0.78
1:AA:336:C:O2'	1:AA:337:C:H5'	1.83	0.78
7:AG:50:ILE:HB	7:AG:58:PRO:HD3	1.64	0.78
27:B1:86:SER:O	27:B1:90:ILE:HD11	1.84	0.78
34:B8:6:THR:CG2	34:B8:63:PRO:HD3	2.14	0.78
35:BA:1290:C:H2'	35:BA:1291:C:H6	1.49	0.78
35:BA:30:G:H2'	35:BA:31:C:H6	1.48	0.78
38:BD:18:VAL:HG23	38:BD:211:ARG:NH2	1.98	0.78
46:BP:17:LYS:HG2	46:BP:17:LYS:O	1.82	0.78
35:BA:2873:A:H1'	48:BR:6:SER:HB2	1.66	0.78
50:BT:91:ARG:HA	50:BT:117:ASP:H	1.49	0.78
4:CD:8:VAL:O	4:CD:10:ARG:N	2.17	0.78
12:CL:70:ILE:N	12:CL:70:ILE:HD12	1.99	0.78
1:CA:658:G:H1'	15:CO:22:THR:HB	1.66	0.78
25:CY:164:ILE:HD12	25:CY:164:ILE:H	1.49	0.78
27:D1:41:ARG:HH11	27:D1:41:ARG:HG3	1.48	0.78
35:DA:1602:U:H3'	35:DA:1603:A:C5'	2.14	0.78
35:DA:2297:C:C2'	35:DA:2298:A:H5'	2.14	0.78
35:DA:2315:G:H2'	35:DA:2316:C:C6	2.18	0.78
26:D0:43:THR:HG22	35:DA:2331:G:O2'	1.84	0.78
38:DD:268:ARG:NH1	38:DD:268:ARG:HB2	1.97	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:31:LYS:NZ	38:DD:31:LYS:HA	1.98	0.78
41:DG:111:LEU:CB	41:DG:112:PRO:HD3	2.10	0.78
43:DI:140:LEU:HD12	43:DI:141:LYS:N	1.98	0.78
46:DP:84:ASN:HA	46:DP:115:LEU:O	1.84	0.78
50:DT:28:VAL:HG13	50:DT:45:PHE:O	1.83	0.78
55:DY:86:ARG:HB3	55:DY:88:LYS:HZ2	1.47	0.78
1:AA:1305:G:H5'	21:AU:4:GLY:HA3	1.64	0.78
1:AA:1442(B):A:N3	1:AA:1442(B):A:H5''	1.99	0.78
2:AB:140:HIS:O	2:AB:143:GLU:HB2	1.84	0.78
4:AD:5:ILE:HG22	4:AD:6:GLY:N	1.99	0.78
35:BA:1019:U:H2'	35:BA:1020:A:C8	2.18	0.78
35:BA:1709:U:H2'	35:BA:1710:C:C6	2.18	0.78
35:BA:598:G:H5'	46:BP:15:ARG:HD2	1.63	0.78
39:BE:108:SER:CB	39:BE:165:VAL:HG21	2.13	0.78
41:BG:170:ARG:NH2	41:BG:182:LYS:HE2	1.99	0.78
46:BP:64:LYS:O	46:BP:66:GLY:N	2.16	0.78
47:BQ:134:ARG:HG2	47:BQ:135:ASP:H	1.49	0.78
54:BX:76:ARG:O	54:BX:76:ARG:HD3	1.83	0.78
4:CD:31:CYS:SG	4:CD:31:CYS:O	2.42	0.78
8:CH:30:ARG:CB	8:CH:30:ARG:HH11	1.95	0.78
39:DE:131:ALA:CB	39:DE:134:ILE:HD11	2.13	0.78
39:DE:100:GLU:O	39:DE:172:VAL:HG23	1.83	0.78
39:DE:52:LEU:HD23	39:DE:75:VAL:CG2	2.13	0.78
51:DU:92:ARG:O	51:DU:94:ASN:N	2.17	0.78
1:AA:1116:C:C2'	1:AA:1117:G:H5''	2.14	0.78
1:AA:1347:G:H22	1:AA:1373:G:H2'	1.47	0.78
1:AA:1497:G:O2'	1:AA:1498:U:H5'	1.84	0.78
13:AM:15:VAL:O	13:AM:19:LEU:HD23	1.84	0.78
22:AV:34:A:H2'	22:AV:35:A:C8	2.19	0.78
35:BA:484:C:H2'	35:BA:485:C:H6	1.46	0.78
37:BC:58:VAL:HG21	37:BC:166:ASP:H	1.47	0.78
35:BA:1245:G:H3'	46:BP:16:ARG:HH22	1.49	0.78
35:BA:1754:C:H5'	50:BT:101:PHE:CD1	2.18	0.78
56:BZ:17:ALA:CA	56:BZ:20:ARG:HB3	2.13	0.78
4:CD:163:GLU:O	4:CD:165:MET:N	2.17	0.78
4:CD:176:LEU:CG	4:CD:177:ASP:H	1.96	0.78
12:CL:47:LYS:CB	12:CL:48:PRO:HD3	2.13	0.78
13:CM:3:ARG:HA	13:CM:9:ILE:HG13	1.64	0.78
35:DA:229:A:H3'	35:DA:230:U:H5'	1.65	0.78
35:DA:2415:G:H4'	46:DP:66:GLY:C	2.04	0.78
35:DA:2691:C:H6	35:DA:2691:C:H5'	1.49	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2312:U:O3'	41:DG:71:THR:HG21	1.83	0.78
44:DN:39:ARG:HD3	44:DN:39:ARG:O	1.84	0.78
48:DR:20:LEU:HD12	48:DR:20:LEU:C	2.05	0.78
11:AK:111:ASP:HA	18:AR:84:LYS:HG3	1.64	0.77
35:BA:1786:A:N7	35:BA:1938:A:N7	2.31	0.77
48:BR:18:LEU:HD13	48:BR:19:ALA:N	1.99	0.77
55:BY:28:LYS:HD2	55:BY:37:VAL:HG12	1.66	0.77
1:CA:405:U:H3'	1:CA:406:G:H5'	1.66	0.77
3:CC:157:ILE:HB	3:CC:164:ARG:NH1	1.99	0.77
11:CK:44:SER:H	11:CK:47:VAL:HG21	1.49	0.77
13:CM:49:THR:HB	13:CM:52:GLU:HG3	1.67	0.77
35:DA:528:A:H2	35:DA:2043:C:C5'	1.96	0.77
38:DD:129:ASN:O	38:DD:193:VAL:HG12	1.84	0.77
39:DE:101:ARG:HD3	39:DE:169:ASN:HD22	1.50	0.77
46:DP:23:PRO:HB2	46:DP:33:ARG:HD3	1.66	0.77
2:AB:200:ILE:HG22	2:AB:202:PRO:HD3	1.66	0.77
5:AE:82:VAL:HG21	5:AE:138:ALA:HA	1.64	0.77
7:AG:113:GLU:HB2	7:AG:119:ARG:HG2	1.65	0.77
10:AJ:4:ILE:HG23	10:AJ:98:ILE:HG23	1.64	0.77
15:AO:54:ARG:O	15:AO:58:MET:HG3	1.84	0.77
35:BA:1230:C:H2'	35:BA:1231:G:C8	2.20	0.77
47:BQ:20:ALA:HA	47:BQ:98:LYS:HB3	1.66	0.77
45:BO:79:PHE:HA	50:BT:72:VAL:HG22	1.65	0.77
55:BY:81:LYS:HG3	55:BY:97:ARG:HG2	1.66	0.77
1:CA:1184:G:H2'	1:CA:1185:G:H8	1.50	0.77
1:CA:1406:U:H2'	1:CA:1407:C:H6	1.49	0.77
4:CD:134:ASP:O	4:CD:136:PRO:HD3	1.83	0.77
12:CL:83:VAL:HG22	12:CL:84:LEU:N	1.98	0.77
16:CP:82:GLN:HE21	16:CP:82:GLN:N	1.82	0.77
25:CY:103:ILE:HD12	25:CY:103:ILE:O	1.84	0.77
28:D2:27:GLU:C	28:D2:29:LYS:H	1.88	0.77
35:DA:1750:G:O2'	35:DA:1751:C:H5'	1.84	0.77
35:DA:196:A:H5''	46:DP:46:LYS:HZ1	1.49	0.77
35:DA:2817:G:H21	35:DA:2836:U:H1'	1.49	0.77
41:DG:144:ILE:CG1	41:DG:145:THR:H	1.96	0.77
52:DV:34:GLU:HB3	52:DV:62:LEU:HD12	1.65	0.77
54:DX:12:VAL:HG11	54:DX:27:THR:HG23	1.66	0.77
5:AE:76:ILE:HG12	5:AE:77:PRO:HD2	1.66	0.77
12:AL:83:VAL:HG22	12:AL:84:LEU:N	1.99	0.77
17:AQ:54:GLY:HA3	17:AQ:82:MET:SD	2.24	0.77
28:B2:56:GLN:HA	28:B2:56:GLN:HE21	1.49	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:B3:56:VAL:HG12	29:B3:57:GLU:N	1.99	0.77
35:BA:1270:C:H5''	35:BA:1271:G:C5'	2.13	0.77
36:BB:74:U:C3'	36:BB:75:G:H5''	2.14	0.77
38:BD:257:LEU:HD23	38:BD:257:LEU:C	2.05	0.77
48:BR:2:ARG:HD2	48:BR:2:ARG:O	1.84	0.77
50:BT:101:PHE:HD2	50:BT:102:ILE:N	1.82	0.77
55:BY:31:LEU:HB2	55:BY:36:ALA:H	1.48	0.77
1:CA:1305:G:H5'	21:CU:4:GLY:HA3	1.65	0.77
12:CL:58:VAL:O	12:CL:65:GLU:HA	1.83	0.77
18:CR:53:ARG:HH12	18:CR:59:SER:HA	1.50	0.77
27:D1:94:LEU:HD13	27:D1:95:LEU:H	1.50	0.77
35:DA:1899:G:N2	35:DA:1902:C:H41	1.82	0.77
35:DA:2715:C:H2'	35:DA:2716:U:C6	2.19	0.77
35:DA:759:G:H2'	35:DA:760:G:H8	1.49	0.77
55:DY:28:LYS:O	55:DY:38:ILE:HB	1.83	0.77
1:AA:1490:C:O2'	1:AA:1491:G:H5'	1.84	0.77
1:AA:686:U:H1'	1:AA:687:A:N7	1.99	0.77
1:AA:806:C:H2'	1:AA:807:A:C8	2.19	0.77
5:AE:19:MET:SD	5:AE:24:ARG:HG2	2.24	0.77
35:BA:2298:A:H2'	35:BA:2299:G:O4'	1.85	0.77
35:BA:268:C:O2	35:BA:268:C:H2'	1.83	0.77
35:BA:2755:C:O2'	35:BA:2756:U:H2'	1.83	0.77
38:BD:133:LEU:HA	38:BD:136:ILE:HD12	1.66	0.77
38:BD:165:ILE:HD12	38:BD:165:ILE:N	1.98	0.77
47:BQ:82:ARG:O	47:BQ:83:MET:HG2	1.84	0.77
49:BS:83:LYS:HG2	49:BS:105:ALA:CB	2.15	0.77
50:BT:109:GLU:HA	50:BT:112:ARG:CG	2.14	0.77
52:BV:5:VAL:HG21	52:BV:36:PRO:HB2	1.66	0.77
1:CA:1424:C:H2'	1:CA:1425:U:C6	2.18	0.77
12:CL:38:THR:HG22	12:CL:57:LYS:O	1.84	0.77
35:DA:1336:A:H2'	35:DA:1337:G:H8	1.50	0.77
35:DA:2716:U:H2'	35:DA:2717:G:C8	2.19	0.77
35:DA:2720:U:O2	35:DA:2720:U:H2'	1.83	0.77
35:DA:729:G:C5	38:DD:208:LYS:HB2	2.20	0.77
48:DR:41:ALA:HB1	48:DR:114:VAL:HG23	1.65	0.77
1:AA:41:G:H2'	1:AA:42:G:C8	2.20	0.77
5:AE:78:HIS:HE1	5:AE:143:ARG:H	1.30	0.77
27:B1:51:VAL:HG23	27:B1:62:VAL:HG11	1.66	0.77
28:B2:52:ASP:O	28:B2:55:ARG:HB3	1.85	0.77
32:B6:15:GLU:OE2	32:B6:41:PRO:HG3	1.85	0.77
38:BD:226:MET:HB3	38:BD:230:ASP:HB2	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:177:ALA:HB1	40:BF:178:PRO:HD2	1.67	0.77
40:BF:63:LYS:HZ1	40:BF:67:GLN:HB3	1.44	0.77
43:BI:102:SER:HB2	43:BI:109:ILE:CG1	2.13	0.77
28:B2:55:ARG:HH22	54:BX:3:THR:HG23	1.49	0.77
6:CF:33:TYR:HD1	6:CF:75:LEU:HG	1.48	0.77
9:CI:21:PRO:HA	9:CI:59:PHE:HA	1.66	0.77
12:CL:119:LYS:HB2	12:CL:120:TYR:CD1	2.19	0.77
20:CT:88:VAL:HA	20:CT:91:LEU:HD12	1.66	0.77
35:DA:1570:A:H2'	35:DA:1571:A:C8	2.19	0.77
35:DA:1722:A:O2'	35:DA:1739:U:H5''	1.84	0.77
35:DA:807:U:H2'	35:DA:808:G:H8	1.49	0.77
46:DP:64:LYS:O	46:DP:66:GLY:N	2.17	0.77
48:DR:99:LYS:O	48:DR:100:LEU:HD22	1.85	0.77
56:DZ:6:LYS:HB2	56:DZ:8:TYR:HE1	1.49	0.77
11:AK:43:SER:HA	11:AK:47:VAL:HG21	1.67	0.77
28:B2:53:LEU:HA	28:B2:56:GLN:CG	2.13	0.77
29:B3:18:ASP:O	29:B3:21:ALA:HB3	1.84	0.77
32:B6:11:LEU:HD11	32:B6:51:GLU:HB2	1.67	0.77
35:BA:1938:A:C2	35:BA:2590:A:H1'	2.20	0.77
38:BD:35:LYS:HE2	38:BD:104:TYR:HB2	1.63	0.77
38:BD:247:ALA:HA	38:BD:254:THR:HG22	1.67	0.77
43:BI:133:HIS:ND1	43:BI:134:PRO:HD2	2.00	0.77
43:BI:140:LEU:HD12	43:BI:141:LYS:N	1.99	0.77
35:BA:666:G:H4'	46:BP:49:ARG:NH2	1.99	0.77
34:B8:25:MET:HE2	46:BP:64:LYS:HG3	1.65	0.77
54:BX:60:ARG:CG	54:BX:72:LYS:H	1.96	0.77
9:CI:50:LEU:HD21	9:CI:81:ILE:CG2	2.14	0.77
18:CR:87:ARG:NH1	18:CR:87:ARG:HB3	1.99	0.77
20:CT:26:ASN:HB2	20:CT:71:THR:HG23	1.66	0.77
31:D5:44:THR:HG22	31:D5:45:VAL:N	2.00	0.77
35:DA:18:C:H2'	35:DA:19:C:H6	1.49	0.77
35:DA:851:U:H2'	35:DA:852:G:H8	1.49	0.77
38:DD:165:ILE:HD12	38:DD:165:ILE:N	1.99	0.77
40:DF:69:HIS:O	40:DF:70:THR:HG23	1.84	0.77
45:DO:119:PRO:HB2	50:DT:68:TYR:CE1	2.19	0.77
46:DP:79:ARG:HH21	46:DP:109:GLY:HA3	1.50	0.77
1:AA:1346:A:H5''	9:AI:120:ARG:HH12	1.49	0.77
4:AD:174:LEU:N	4:AD:186:LEU:HD12	2.00	0.77
9:AI:95:LYS:HD3	9:AI:96:LEU:N	1.98	0.77
11:AK:121:PRO:O	11:AK:126:ARG:HB2	1.85	0.77
26:B0:43:THR:HG22	35:BA:2331:G:O2'	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:8:ASN:HD22	33:B7:9:ARG:N	1.83	0.77
35:BA:467:G:H2'	35:BA:468:G:H8	1.50	0.77
35:BA:925:C:H2'	35:BA:926:A:C5'	2.14	0.77
38:BD:163:ALA:HB1	38:BD:175:LEU:HD21	1.65	0.77
41:BG:114:ILE:HD12	41:BG:117:PHE:CG	2.19	0.77
41:BG:73:ALA:HB2	41:BG:87:PRO:HG2	1.65	0.77
45:BO:2:ILE:HD11	45:BO:82:ASN:HB3	1.66	0.77
34:B8:13:ARG:CB	46:BP:63:PRO:HA	2.15	0.77
1:CA:1190:G:P	3:CC:5:ILE:HG23	2.25	0.77
11:CK:21:ILE:HD12	11:CK:21:ILE:N	1.99	0.77
11:CK:17:GLY:HA3	11:CK:80:VAL:HA	1.65	0.77
25:CY:52:LEU:HD21	25:CY:56:ALA:HB3	1.65	0.77
34:D8:23:VAL:HG13	34:D8:47:LYS:O	1.84	0.77
34:D8:32:LEU:O	34:D8:33:ASN:HB3	1.84	0.77
35:DA:1270:C:H5''	35:DA:1271:G:C5'	2.14	0.77
38:DD:167:GLY:H	38:DD:174:ILE:HB	1.50	0.77
41:DG:102:PHE:CE1	41:DG:106:LEU:HD13	2.19	0.77
43:DI:79:ILE:CG1	43:DI:140:LEU:HD11	2.14	0.77
50:DT:96:ARG:HG2	50:DT:96:ARG:HH11	1.49	0.77
52:DV:15:GLU:HB3	52:DV:16:PRO:HD2	1.66	0.77
1:AA:405:U:H3'	1:AA:406:G:H5'	1.67	0.77
1:AA:665:A:H2'	1:AA:725:G:N2	2.00	0.77
4:AD:59:ARG:HH22	4:AD:66:ARG:HH22	1.31	0.77
1:AA:528:C:H41	12:AL:49:ASN:HD22	1.29	0.77
35:BA:2521:C:H42	35:BA:2544:G:H1	1.32	0.77
35:BA:491:G:H2'	35:BA:492:A:C8	2.19	0.77
39:BE:33:VAL:CG1	39:BE:89:ASP:H	1.95	0.77
41:BG:47:LYS:HE3	41:BG:81:LYS:HB3	1.66	0.77
46:BP:79:ARG:HH21	46:BP:109:GLY:HA3	1.50	0.77
1:CA:662:G:H2'	1:CA:663:A:H8	1.49	0.77
3:CC:6:HIS:ND1	14:CN:49:HIS:HB3	2.00	0.77
9:CI:9:ARG:HG3	9:CI:14:VAL:HG22	1.67	0.77
16:CP:45:THR:HG22	16:CP:47:ASP:H	1.50	0.77
35:DA:8:A:H2'	35:DA:9:U:H6	1.50	0.77
38:DD:231:HIS:ND1	38:DD:232:PRO:HD2	1.99	0.77
35:DA:626:U:O2	46:DP:105:LEU:HG	1.84	0.77
6:AF:76:ALA:HB1	6:AF:80:ARG:NH2	1.99	0.77
8:AH:25:ASP:HA	8:AH:59:LEU:O	1.84	0.77
10:AJ:4:ILE:HB	10:AJ:74:ILE:CD1	2.15	0.77
11:AK:59:TYR:CZ	11:AK:63:LEU:HD11	2.20	0.77
19:AS:70:LYS:HB3	19:AS:70:LYS:NZ	1.99	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:192:U:H4'	20:AT:103:GLY:HA2	1.65	0.77
24:AX:13:A:O2'	24:AX:14:U:H5'	1.85	0.77
31:B5:25:LEU:HD11	53:BW:19:LEU:HB3	1.66	0.77
35:BA:1635:G:H8	35:BA:1635:G:H5'	1.48	0.77
35:BA:2820:A:C4'	48:BR:5:LYS:HE2	2.15	0.77
35:BA:2854:G:H2'	35:BA:2855:C:C6	2.19	0.77
35:BA:963:U:H2'	35:BA:964:C:C6	2.20	0.77
41:BG:115:ARG:HH22	41:BG:136:ARG:HD2	1.48	0.77
47:BQ:39:PRO:HB3	47:BQ:99:PRO:CD	2.10	0.77
56:BZ:110:GLY:C	56:BZ:112:ARG:N	2.37	0.77
1:CA:1478:C:H2'	1:CA:1479:C:C6	2.20	0.77
9:CI:63:ILE:HD12	9:CI:63:ILE:H	1.50	0.77
28:D2:30:ARG:H	28:D2:30:ARG:HD2	1.48	0.77
31:D5:57:VAL:HG23	31:D5:58:LEU:H	1.50	0.77
35:DA:1378:A:H4'	35:DA:1379:A:OP1	1.82	0.77
35:DA:2558:C:H2'	35:DA:2559:C:C6	2.20	0.77
39:DE:37:ARG:HB2	39:DE:46:ALA:HB3	1.66	0.77
41:DG:41:GLN:NE2	41:DG:153:ARG:HB3	2.00	0.77
1:AA:59:A:C5'	1:AA:60:A:H5''	2.15	0.77
2:AB:36:ARG:HB3	2:AB:41:ILE:HD11	1.67	0.77
7:AG:50:ILE:O	7:AG:54:THR:HG23	1.84	0.77
9:AI:65:VAL:O	9:AI:66:ARG:HG3	1.85	0.77
35:BA:581:C:H2'	35:BA:582:G:H8	1.50	0.77
39:BE:52:LEU:HD23	39:BE:75:VAL:CG2	2.15	0.77
40:BF:132:VAL:HG22	40:BF:133:ASN:N	2.00	0.77
44:BN:43:THR:HB	44:BN:46:VAL:HB	1.67	0.77
4:CD:30:LYS:C	4:CD:32:ALA:N	2.39	0.77
20:CT:104:LEU:HD23	20:CT:105:SER:N	1.99	0.77
34:D8:32:LEU:HG	34:D8:34:TRP:HE3	1.50	0.77
35:DA:1635:G:H8	35:DA:1635:G:H5'	1.50	0.77
44:DN:74:ARG:CZ	44:DN:101:HIS:HB3	2.15	0.77
49:DS:27:SER:HA	49:DS:89:ARG:HD2	1.66	0.77
50:DT:23:ARG:HG2	50:DT:120:ARG:NH1	2.00	0.77
51:DU:88:ILE:C	51:DU:90:VAL:N	2.38	0.77
1:AA:1410:G:H2'	1:AA:1411:C:H6	1.50	0.76
2:AB:95:GLN:HG3	2:AB:148:TYR:HA	1.66	0.76
9:AI:63:ILE:HD12	9:AI:63:ILE:H	1.48	0.76
35:BA:1364:G:H1'	35:BA:1368:G:N2	1.99	0.76
35:BA:1401:G:H2'	35:BA:1402:C:C6	2.20	0.76
39:BE:6:GLY:HA2	39:BE:51:PHE:CE2	2.19	0.76
49:BS:26:LEU:O	49:BS:26:LEU:HD23	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1298:C:H4'	1:CA:1299:A:O4'	1.86	0.76
1:CA:198:G:H2'	1:CA:199:G:C8	2.20	0.76
1:CA:553:A:H2'	1:CA:554:C:C6	2.20	0.76
8:CH:53:VAL:O	8:CH:56:LYS:HB2	1.84	0.76
18:CR:56:THR:HB	18:CR:58:LEU:CD1	2.13	0.76
28:D2:17:SER:OG	28:D2:18:PRO:HD3	1.83	0.76
38:DD:172:TYR:HD1	38:DD:186:HIS:HA	1.50	0.76
38:DD:27:THR:O	38:DD:28:GLU:HB2	1.83	0.76
35:DA:2562:U:H1'	45:DO:23:ARG:NH1	1.99	0.76
46:DP:112:LEU:HD22	46:DP:113:LYS:H	1.50	0.76
56:DZ:120:ILE:O	56:DZ:121:HIS:HB2	1.83	0.76
1:AA:1406:U:H2'	1:AA:1407:C:O4'	1.86	0.76
18:AR:52:PRO:O	18:AR:56:THR:HG23	1.85	0.76
25:AY:140:LEU:HD11	25:AY:157:ALA:HB1	1.65	0.76
27:B1:26:ARG:HB2	27:B1:34:THR:OG1	1.85	0.76
35:BA:2415:G:H4'	46:BP:66:GLY:C	2.06	0.76
35:BA:2730:C:O2'	35:BA:2731:G:H5'	1.85	0.76
35:BA:2779:U:H1'	35:BA:2781:A:C5	2.20	0.76
35:BA:587:C:C5	46:BP:33:ARG:HD2	2.20	0.76
35:BA:705:A:H1'	38:BD:9:TYR:CE1	2.19	0.76
38:BD:44:ASN:CB	38:BD:49:ILE:HA	2.15	0.76
39:BE:101:ARG:HD3	39:BE:169:ASN:HD22	1.49	0.76
39:BE:37:ARG:HB2	39:BE:46:ALA:HB3	1.67	0.76
42:BH:67:LEU:HG	42:BH:71:LEU:HD21	1.67	0.76
43:BI:77:LEU:HB2	43:BI:140:LEU:CD1	2.13	0.76
45:BO:6:THR:HG22	45:BO:7:TYR:H	1.50	0.76
56:BZ:103:ARG:HG3	56:BZ:136:PHE:CZ	2.20	0.76
1:CA:1073:U:H2'	1:CA:1074:G:H8	1.49	0.76
1:CA:778:G:O2'	1:CA:779:C:H5'	1.84	0.76
2:CB:29:ALA:O	2:CB:31:TYR:N	2.18	0.76
9:CI:95:LYS:HD3	9:CI:96:LEU:N	2.01	0.76
1:CA:255:G:H1'	17:CQ:16:GLN:NE2	1.99	0.76
35:DA:1314:C:H6	35:DA:1314:C:H5'	1.51	0.76
35:DA:1639:U:H2'	35:DA:1640:C:H5''	1.66	0.76
35:DA:1717:G:H3'	35:DA:1718:G:H5''	1.66	0.76
35:DA:2036:C:C6	35:DA:2036:C:H5'	2.16	0.76
31:D5:2:ALA:HB3	35:DA:747:U:C2	2.20	0.76
35:DA:2591:C:OP2	38:DD:239:ARG:HB2	1.85	0.76
41:DG:141:PHE:CD1	41:DG:142:PRO:HD2	2.20	0.76
1:AA:1074:G:H2'	1:AA:1075:C:H6	1.50	0.76
9:AI:112:LYS:HA	9:AI:119:ALA:HB2	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AT:100:ILE:HD12	20:AT:100:ILE:N	1.99	0.76
35:BA:1722:A:O2'	35:BA:1739:U:H5''	1.84	0.76
35:BA:2052:G:H2'	35:BA:2053:G:H8	1.51	0.76
35:BA:2220:G:H2'	35:BA:2221:G:H8	1.48	0.76
39:BE:167:VAL:HG22	39:BE:168:MET:H	1.50	0.76
40:BF:53:THR:CG2	40:BF:56:GLU:HB2	2.15	0.76
51:BU:92:ARG:O	51:BU:94:ASN:N	2.18	0.76
47:BQ:108:GLY:HA3	56:BZ:116:VAL:HG21	1.67	0.76
1:CA:313:A:H2'	1:CA:314:C:C6	2.20	0.76
1:CA:764:C:H2'	1:CA:765:G:H8	1.50	0.76
2:CB:19:HIS:HA	2:CB:39:ILE:HD13	1.67	0.76
2:CB:36:ARG:HB3	2:CB:41:ILE:HD11	1.66	0.76
2:CB:71:VAL:HG21	2:CB:93:VAL:HG23	1.66	0.76
35:DA:1112:G:H1'	35:DA:1113:U:OP1	1.85	0.76
35:DA:1639:U:C2'	35:DA:1640:C:H5''	2.15	0.76
35:DA:1914:C:H2'	35:DA:1915:U:O4'	1.85	0.76
35:DA:97:C:H2'	35:DA:98:G:C8	2.21	0.76
38:DD:257:LEU:HD23	38:DD:257:LEU:C	2.06	0.76
34:D8:25:MET:HE2	46:DP:64:LYS:HG3	1.67	0.76
1:AA:1088:G:H2'	1:AA:1089:G:H8	1.50	0.76
1:AA:1423:G:H5'	45:BO:49:ARG:NH2	2.01	0.76
8:AH:86:ILE:HG12	8:AH:135:CYS:HA	1.65	0.76
9:AI:50:LEU:HD21	9:AI:81:ILE:CG2	2.14	0.76
35:BA:8:A:H2'	35:BA:9:U:H6	1.50	0.76
46:BP:83:VAL:HG12	46:BP:112:LEU:HD21	1.66	0.76
51:BU:88:ILE:C	51:BU:90:VAL:H	1.87	0.76
54:BX:57:LEU:HD12	54:BX:76:ARG:NE	2.00	0.76
1:CA:1435:G:H2'	1:CA:1436:U:C5	2.20	0.76
1:CA:735:C:O2'	1:CA:736:C:H5'	1.85	0.76
9:CI:112:LYS:HA	9:CI:119:ALA:HB2	1.67	0.76
1:CA:528:C:H41	12:CL:49:ASN:HD22	1.34	0.76
35:DA:1709:U:H2'	35:DA:1710:C:H6	1.49	0.76
31:D5:2:ALA:HA	35:DA:2015:A:C1'	2.15	0.76
35:DA:2755:C:O2'	35:DA:2756:U:H2'	1.86	0.76
35:DA:813:U:H2'	35:DA:814:C:C6	2.21	0.76
40:DF:3:GLU:HB2	40:DF:24:LEU:HG	1.68	0.76
47:DQ:82:ARG:O	47:DQ:83:MET:HG2	1.85	0.76
51:DU:18:LEU:O	51:DU:18:LEU:HD23	1.86	0.76
51:DU:6:THR:HG21	51:DU:10:ARG:HH21	1.48	0.76
35:DA:1614:A:C2	53:DW:87:PRO:HB3	2.20	0.76
1:AA:1298:C:H4'	1:AA:1299:A:O4'	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:19:GLN:HE21	35:BA:379:G:N2	1.81	0.76
35:BA:742:G:H2'	35:BA:743:G:C8	2.18	0.76
39:BE:30:PRO:HD3	39:BE:180:ASN:ND2	1.99	0.76
45:BO:63:VAL:CG2	45:BO:84:ALA:HA	2.13	0.76
48:BR:103:ARG:HG2	48:BR:103:ARG:HH11	1.50	0.76
54:BX:36:LYS:O	54:BX:38:GLU:N	2.18	0.76
1:CA:1053:G:O6	1:CA:1200:C:H5''	1.86	0.76
1:CA:646:U:H2'	1:CA:647:C:C6	2.20	0.76
3:CC:86:VAL:O	3:CC:90:GLU:HG2	1.86	0.76
8:CH:109:ILE:HG12	8:CH:110:ALA:N	2.00	0.76
16:CP:14:ASN:N	16:CP:15:PRO:HD3	1.99	0.76
17:CQ:97:SER:O	17:CQ:98:LEU:HG	1.84	0.76
27:D1:47:GLN:HB3	27:D1:64:ALA:HB2	1.65	0.76
35:DA:2243:U:H2'	35:DA:2244:U:H6	1.51	0.76
35:DA:431:U:O2'	35:DA:432:A:H5'	1.85	0.76
35:DA:558:G:H2'	35:DA:559:G:H8	1.50	0.76
35:DA:877:U:H2'	35:DA:878:A:H5''	1.65	0.76
38:DD:65:ILE:HD12	38:DD:65:ILE:C	2.06	0.76
40:DF:185:ASP:HA	40:DF:188:ARG:HB3	1.67	0.76
35:DA:2415:G:H4'	46:DP:66:GLY:CA	2.15	0.76
50:DT:65:LYS:HA	50:DT:65:LYS:HZ2	1.48	0.76
56:DZ:33:LEU:HD21	56:DZ:35:ARG:HB2	1.65	0.76
1:AA:1239:A:H62	1:AA:1299:A:H62	1.33	0.76
1:AA:1366:C:H2'	1:AA:1367:C:C6	2.21	0.76
3:AC:6:HIS:ND1	14:AN:49:HIS:HB3	2.00	0.76
8:AH:120:THR:OG1	8:AH:123:GLU:HG3	1.86	0.76
10:AJ:48:THR:HG22	10:AJ:49:VAL:H	1.51	0.76
25:AY:61:PRO:HD2	25:AY:65:THR:O	1.86	0.76
31:B5:57:VAL:HG23	31:B5:58:LEU:H	1.49	0.76
33:B7:34:ARG:NH1	33:B7:42:LEU:O	2.17	0.76
34:B8:39:LYS:HG2	34:B8:42:ARG:NH1	2.00	0.76
34:B8:23:VAL:HG12	34:B8:46:ARG:HH12	1.51	0.76
38:BD:34:VAL:HG22	38:BD:35:LYS:HG3	1.66	0.76
42:BH:85:LYS:O	42:BH:132:ARG:HA	1.85	0.76
36:BB:7:G:H4'	49:BS:29:PHE:CD2	2.20	0.76
1:CA:1116:C:C2'	1:CA:1117:G:H5''	2.16	0.76
1:CA:1299:A:H2'	1:CA:1301:U:C6	2.20	0.76
1:CA:976:G:N2	1:CA:1362:C:H2'	2.01	0.76
1:CA:636:U:H2'	1:CA:637:G:C8	2.20	0.76
2:CB:95:GLN:HG3	2:CB:148:TYR:HA	1.67	0.76
8:CH:25:ASP:HA	8:CH:59:LEU:O	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:873:G:H2'	35:DA:874:G:C8	2.20	0.76
38:DD:48:ARG:HH11	38:DD:48:ARG:HG3	1.51	0.76
39:DE:102:VAL:HA	39:DE:200:GLU:HA	1.68	0.76
34:D8:13:ARG:HB3	46:DP:63:PRO:CA	2.14	0.76
1:AA:198:G:H2'	1:AA:199:G:C8	2.20	0.76
2:AB:19:HIS:HA	2:AB:39:ILE:HD13	1.68	0.76
1:CA:328:C:H4'	1:CA:329:A:H5'	1.67	0.76
1:CA:543:C:H2'	1:CA:544:G:H8	1.50	0.76
12:CL:54:LYS:N	12:CL:54:LYS:HD2	2.00	0.76
18:CR:86:VAL:HG12	18:CR:87:ARG:HH12	1.50	0.76
24:CX:16:U:H2'	24:CX:17:U:C6	2.20	0.76
35:DA:259:G:H21	35:DA:621:A:H8	1.33	0.76
35:DA:2730:C:O2'	35:DA:2731:G:H5'	1.84	0.76
35:DA:588:U:H6	35:DA:588:U:O5'	1.68	0.76
42:DH:85:LYS:O	42:DH:132:ARG:HA	1.86	0.76
54:DX:36:LYS:O	54:DX:38:GLU:N	2.18	0.76
1:AA:255:G:H1'	17:AQ:16:GLN:NE2	1.99	0.76
35:BA:2201:C:O2'	35:BA:2202:C:H5'	1.85	0.76
44:BN:58:ASP:C	44:BN:60:ILE:H	1.88	0.76
46:BP:59:LEU:HA	46:BP:61:ARG:CZ	2.16	0.76
47:BQ:119:ARG:HG2	47:BQ:120:ILE:HD13	1.66	0.76
49:BS:89:ARG:HA	49:BS:89:ARG:NE	2.01	0.76
50:BT:23:ARG:HG2	50:BT:120:ARG:NH1	2.01	0.76
2:CB:72:GLY:HA3	2:CB:165:VAL:HG13	1.68	0.76
4:CD:5:ILE:O	4:CD:6:GLY:O	2.04	0.76
40:DF:177:ALA:HB1	40:DF:178:PRO:HD2	1.67	0.76
41:DG:39:ILE:HA	41:DG:157:ILE:CA	2.16	0.76
42:DH:105:LEU:HD21	42:DH:113:VAL:HB	1.68	0.76
46:DP:38:GLN:CG	46:DP:39:LYS:H	1.99	0.76
47:DQ:34:LEU:HD12	47:DQ:35:VAL:H	1.50	0.76
50:DT:109:GLU:HA	50:DT:112:ARG:CG	2.15	0.76
50:DT:91:ARG:HA	50:DT:117:ASP:H	1.50	0.76
1:AA:854:G:H3'	1:AA:871:U:O4	1.86	0.76
3:AC:9:GLY:O	3:AC:12:LEU:HB2	1.85	0.76
12:AL:84:LEU:HD23	12:AL:85:ILE:N	2.01	0.76
25:AY:170:ALA:O	25:AY:174:GLN:HB3	1.86	0.76
25:AY:173:ASP:O	25:AY:177:GLU:HB2	1.85	0.76
27:B1:33:LYS:HG2	27:B1:34:THR:H	1.50	0.76
34:B8:32:LEU:HG	34:B8:34:TRP:HE3	1.50	0.76
35:BA:1639:U:H2'	35:BA:1640:C:H5''	1.67	0.76
35:BA:2821:A:H2'	35:BA:2822:G:C8	2.20	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:528:A:N1	35:BA:2042:A:H2'	2.01	0.76
35:BA:564:C:H2'	35:BA:565:C:C6	2.21	0.76
39:BE:167:VAL:O	39:BE:168:MET:HG3	1.84	0.76
39:BE:98:PRO:HG3	39:BE:175:VAL:HG12	1.67	0.76
43:BI:79:ILE:CG1	43:BI:140:LEU:HD11	2.16	0.76
44:BN:34:LEU:CD2	44:BN:120:LEU:HB2	2.14	0.76
1:CA:600:C:O2'	1:CA:601:C:H5'	1.85	0.76
6:CF:53:ALA:HB3	6:CF:86:ARG:HH12	1.48	0.76
17:CQ:40:LYS:HD2	17:CQ:42:TYR:CE1	2.20	0.76
25:CY:70:SER:HB3	25:CY:76:LEU:HB2	1.66	0.76
35:DA:2762:G:H2'	35:DA:2763:G:H5''	1.68	0.76
38:DD:135:PHE:HD1	38:DD:135:PHE:H	1.33	0.76
38:DD:45:ASN:ND2	38:DD:46:GLN:H	1.84	0.76
38:DD:45:ASN:CG	38:DD:46:GLN:H	1.85	0.76
35:DA:2811:G:OP1	39:DE:60:ASN:HB2	1.85	0.76
44:DN:34:LEU:HD21	44:DN:120:LEU:CB	2.15	0.76
54:DX:60:ARG:CG	54:DX:72:LYS:H	1.99	0.76
55:DY:28:LYS:HZ2	55:DY:37:VAL:HG12	1.49	0.76
1:AA:1489:G:H2'	1:AA:1490:C:H6	1.51	0.76
7:AG:79:ARG:HE	7:AG:84:ASN:ND2	1.82	0.76
12:AL:58:VAL:O	12:AL:65:GLU:HA	1.84	0.76
25:AY:70:SER:HB3	25:AY:76:LEU:CD1	2.16	0.76
31:B5:16:ARG:HG2	31:B5:16:ARG:HH11	1.50	0.76
35:BA:769:G:O2'	35:BA:770:G:H5'	1.86	0.76
35:BA:873:G:H2'	35:BA:874:G:C8	2.21	0.76
45:BO:105:GLU:OE1	45:BO:105:GLU:N	2.19	0.76
46:BP:112:LEU:HD22	46:BP:113:LYS:H	1.51	0.76
35:BA:2415:G:H4'	46:BP:66:GLY:CA	2.15	0.76
51:BU:93:LYS:H	51:BU:93:LYS:HD3	1.50	0.76
54:BX:64:LYS:HG2	54:BX:65:ARG:N	1.99	0.76
56:BZ:28:MET:CE	56:BZ:37:VAL:HG21	2.14	0.76
56:BZ:71:VAL:HG22	56:BZ:88:PHE:HE2	1.49	0.76
1:CA:348:G:O2'	1:CA:349:A:H5'	1.86	0.76
1:CA:59:A:C5'	1:CA:60:A:H5''	2.15	0.76
2:CB:169:LYS:HD3	2:CB:169:LYS:C	2.06	0.76
2:CB:200:ILE:HG22	2:CB:202:PRO:HD3	1.66	0.76
19:CS:70:LYS:NZ	19:CS:70:LYS:HB3	2.00	0.76
20:CT:100:ILE:HD12	20:CT:100:ILE:N	2.00	0.76
20:CT:38:LYS:O	20:CT:41:ILE:HG12	1.86	0.76
35:DA:1364:G:H1'	35:DA:1368:G:N2	2.01	0.76
35:DA:740:U:H2'	35:DA:741:G:H8	1.45	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:860:U:H5	35:DA:917:A:N7	1.84	0.76
38:DD:35:LYS:HE2	38:DD:104:TYR:HB2	1.68	0.76
41:DG:37:VAL:CG1	41:DG:94:LEU:HD12	2.16	0.76
47:DQ:134:ARG:HG2	47:DQ:135:ASP:H	1.51	0.76
39:DE:183:LEU:HD21	50:DT:11:GLU:HB3	1.68	0.76
56:DZ:169:GLU:HG2	56:DZ:170:THR:N	2.00	0.76
1:AA:1271:G:H5'	1:AA:1314:C:H5''	1.67	0.75
1:AA:770:C:O2'	1:AA:771:G:H5'	1.86	0.75
4:AD:5:ILE:HG22	4:AD:6:GLY:H	1.51	0.75
9:AI:21:PRO:HA	9:AI:59:PHE:HA	1.66	0.75
27:B1:94:LEU:HD13	27:B1:95:LEU:O	1.85	0.75
35:BA:1176:G:H1'	35:BA:1177:A:OP1	1.86	0.75
35:BA:2487:G:H2'	35:BA:2488:A:H8	1.51	0.75
38:BD:27:THR:O	38:BD:28:GLU:HB2	1.86	0.75
40:BF:185:ASP:HA	40:BF:188:ARG:HB3	1.67	0.75
43:BI:98:ALA:HB1	43:BI:109:ILE:CB	2.15	0.75
56:BZ:125:LEU:HD23	56:BZ:126:VAL:N	2.00	0.75
7:CG:113:GLU:HB2	7:CG:119:ARG:HG2	1.68	0.75
13:CM:23:TYR:CE1	13:CM:71:ARG:HB2	2.21	0.75
18:CR:50:ILE:HD12	18:CR:70:ILE:HG21	1.67	0.75
31:D5:15:ARG:O	31:D5:18:ALA:HB3	1.86	0.75
31:D5:45:VAL:HG22	31:D5:51:TYR:HD1	1.49	0.75
35:DA:2873:A:H1'	48:DR:6:SER:HB2	1.66	0.75
35:DA:30:G:H2'	35:DA:31:C:H6	1.50	0.75
35:DA:925:C:H2'	35:DA:926:A:C5'	2.16	0.75
39:DE:181:LEU:HD22	39:DE:181:LEU:N	2.01	0.75
41:DG:34:LEU:HD12	41:DG:99:MET:SD	2.25	0.75
41:DG:46:ALA:HB2	41:DG:88:ILE:HB	1.67	0.75
45:DO:121:VAL:C	45:DO:122:LEU:HD12	2.06	0.75
1:AA:625:G:H2'	1:AA:626:U:C6	2.20	0.75
1:AA:878:G:H5'	8:AH:89:PRO:HG2	1.68	0.75
12:AL:54:LYS:HD2	12:AL:54:LYS:N	2.00	0.75
28:B2:33:MET:HG2	54:BX:10:ALA:HB2	1.68	0.75
32:B6:11:LEU:HG	32:B6:26:ASN:ND2	2.00	0.75
35:BA:1270:C:H5''	35:BA:1271:G:O5'	1.85	0.75
35:BA:1484:G:C3'	35:BA:1485:G:H5''	2.16	0.75
35:BA:154:G:H1	35:BA:172:C:H42	1.34	0.75
35:BA:1750:G:O2'	35:BA:1751:C:H5'	1.86	0.75
35:BA:18:C:H2'	35:BA:19:C:H6	1.51	0.75
35:BA:2720:U:H2'	35:BA:2720:U:O2	1.84	0.75
35:BA:877:U:H2'	35:BA:878:A:H5''	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BB:91:C:H2'	36:BB:92:C:H6	1.51	0.75
38:BD:65:ILE:HD12	38:BD:65:ILE:C	2.07	0.75
1:CA:429:U:H1'	1:CA:430:A:H5''	1.69	0.75
1:CA:551:U:H2'	1:CA:552:U:C6	2.21	0.75
35:DA:1509(B):A:H2'	35:DA:1510:G:H8	1.50	0.75
35:DA:1578:U:C3'	35:DA:1579:A:H5''	2.16	0.75
35:DA:2103:C:C3'	35:DA:2104:G:H5''	2.16	0.75
35:DA:2298:A:H2'	35:DA:2299:G:O4'	1.86	0.75
35:DA:2502:G:H5'	35:DA:2503:A:H5''	1.68	0.75
35:DA:904:C:H2'	35:DA:904:C:O2	1.85	0.75
46:DP:47:ASP:HB3	46:DP:48:PRO:O	1.85	0.75
1:AA:403:C:H2'	1:AA:404:U:H6	1.52	0.75
8:AH:11:THR:HA	8:AH:14:ARG:NH1	2.01	0.75
25:AY:65:THR:HG23	25:AY:101:ILE:O	1.86	0.75
27:B1:32:LYS:HA	35:BA:2396:G:HO2'	1.52	0.75
35:BA:1446:C:H42	35:BA:1465:G:H1	1.34	0.75
35:BA:2243:U:H2'	35:BA:2244:U:H6	1.50	0.75
39:BE:128:SER:O	39:BE:129:HIS:HB2	1.86	0.75
41:BG:93:THR:HG22	41:BG:95:ARG:HG3	1.68	0.75
44:BN:34:LEU:HD21	44:BN:120:LEU:CB	2.13	0.75
54:BX:51:VAL:HG13	54:BX:81:VAL:H	1.50	0.75
55:BY:79:CYS:SG	55:BY:80:GLY:N	2.59	0.75
1:CA:385:C:O2'	1:CA:386:C:H5'	1.86	0.75
1:CA:522:C:N4	1:CA:528:C:H42	1.84	0.75
12:CL:89:ARG:HB2	12:CL:89:ARG:NH1	2.00	0.75
23:CW:27:G:H2'	23:CW:28:U:C6	2.21	0.75
35:DA:2303:G:H4'	41:DG:124:SER:O	1.85	0.75
35:DA:2857:G:N2	35:DA:2859:G:H3'	2.01	0.75
35:DA:536:A:H2'	35:DA:537:C:C6	2.21	0.75
35:DA:792:G:H5''	35:DA:793:A:H5'	1.69	0.75
41:DG:110:ALA:CA	41:DG:140:ILE:HD13	2.16	0.75
44:DN:23:LEU:HA	44:DN:26:LEU:HB3	1.66	0.75
47:DQ:16:ARG:HB3	47:DQ:16:ARG:NH1	2.01	0.75
56:DZ:79:ARG:HG2	56:DZ:80:ARG:HG2	1.68	0.75
1:AA:16:A:O2'	1:AA:17:U:H5'	1.87	0.75
1:AA:337:C:H2'	1:AA:338:A:H8	1.50	0.75
1:AA:452:A:H4'	16:AP:72:ARG:NH2	2.02	0.75
20:AT:88:VAL:HA	20:AT:91:LEU:HD12	1.65	0.75
35:BA:2297:C:C2'	35:BA:2298:A:H5'	2.16	0.75
35:BA:2307:G:N2	35:BA:2308:G:H5'	2.01	0.75
35:BA:2533:A:C2'	35:BA:2534:A:H5''	2.16	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:131:TYR:OH	41:BG:133:LEU:HD23	1.85	0.75
1:CA:1088:G:H2'	1:CA:1089:G:H8	1.49	0.75
1:CA:375:U:H2'	1:CA:376:G:H8	1.52	0.75
1:CA:428:G:H4'	1:CA:429:U:O5'	1.87	0.75
5:CE:145:LYS:O	5:CE:148:VAL:HB	1.87	0.75
6:CF:10:LEU:HA	6:CF:84:ASN:O	1.86	0.75
33:D7:34:ARG:NH1	33:D7:42:LEU:O	2.20	0.75
35:DA:1403:C:H5''	35:DA:1471:A:C1'	2.16	0.75
35:DA:2781:A:H5''	35:DA:2782:G:H5'	1.68	0.75
41:DG:105:LYS:HB2	41:DG:105:LYS:HZ2	1.52	0.75
54:DX:5:TYR:O	54:DX:7:VAL:N	2.19	0.75
3:AC:73:PRO:O	3:AC:76:VAL:HG22	1.86	0.75
13:AM:49:THR:HB	13:AM:52:GLU:HG3	1.67	0.75
25:AY:76:LEU:HA	25:AY:79:ILE:HD12	1.68	0.75
31:B5:42:PRO:HB2	35:BA:2815:C:O2'	1.86	0.75
35:BA:2729:G:H1'	39:BE:187:ALA:CB	2.16	0.75
40:BF:114:VAL:HG23	40:BF:115:ALA:N	2.01	0.75
55:BY:28:LYS:HZ2	55:BY:37:VAL:HG12	1.51	0.75
1:CA:1347:G:H22	1:CA:1373:G:H2'	1.47	0.75
1:CA:1436:U:H2'	1:CA:1437:C:C6	2.21	0.75
1:CA:735:C:H2'	1:CA:736:C:H6	1.51	0.75
9:CI:65:VAL:O	9:CI:66:ARG:HG3	1.87	0.75
13:CM:108:ARG:N	13:CM:108:ARG:HD2	2.02	0.75
13:CM:90:LEU:HA	13:CM:93:ARG:HB2	1.68	0.75
16:CP:26:ARG:NH1	16:CP:26:ARG:HB3	2.01	0.75
19:CS:36:ARG:NH2	19:CS:75:ALA:HB3	2.00	0.75
25:CY:68:VAL:CG2	25:CY:99:LEU:HB2	2.16	0.75
31:D5:16:ARG:HH12	31:D5:17:ASP:CG	1.89	0.75
34:D8:23:VAL:HG12	34:D8:46:ARG:NH1	2.00	0.75
35:DA:1938:A:C2	35:DA:2590:A:H1'	2.21	0.75
38:DD:226:MET:HB3	38:DD:230:ASP:HB2	1.67	0.75
51:DU:68:ALA:O	51:DU:71:GLN:HB3	1.87	0.75
55:DY:81:LYS:HG3	55:DY:97:ARG:HG2	1.67	0.75
1:AA:348:G:O2'	1:AA:349:A:H5'	1.85	0.75
6:AF:48:LEU:HD22	18:AR:77:GLY:HA3	1.69	0.75
6:AF:50:TYR:CE1	18:AR:77:GLY:HA2	2.22	0.75
13:AM:23:TYR:CE1	13:AM:71:ARG:HB2	2.21	0.75
14:AN:16:PHE:CD2	14:AN:16:PHE:N	2.53	0.75
25:AY:7:TYR:CE2	25:AY:160:GLU:HG2	2.22	0.75
35:BA:1112:G:H1'	35:BA:1113:U:OP1	1.86	0.75
35:BA:2716:U:H2'	35:BA:2717:G:C8	2.20	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:170:LEU:HD12	40:BF:171:PRO:HD2	1.69	0.75
49:BS:85:VAL:HG23	49:BS:86:ALA:H	1.52	0.75
52:BV:70:ILE:CB	52:BV:90:PRO:HB2	2.17	0.75
55:BY:75:ILE:HD13	55:BY:76:CYS:N	2.01	0.75
4:CD:174:LEU:N	4:CD:186:LEU:HD12	2.02	0.75
29:D3:18:ASP:O	29:D3:21:ALA:HB3	1.87	0.75
29:D3:59:VAL:HG12	29:D3:60:GLU:N	2.02	0.75
35:DA:1709:U:H2'	35:DA:1710:C:C6	2.21	0.75
35:DA:1937:A:O2'	35:DA:1938:A:H5'	1.87	0.75
40:DF:53:THR:H	40:DF:56:GLU:CB	1.99	0.75
49:DS:13:ARG:N	49:DS:13:ARG:HD2	2.02	0.75
49:DS:83:LYS:HG2	49:DS:105:ALA:CB	2.15	0.75
50:DT:101:PHE:HD2	50:DT:102:ILE:N	1.83	0.75
52:DV:61:VAL:HB	52:DV:99:ILE:N	2.01	0.75
1:AA:1116:C:H2'	1:AA:1117:G:C5'	2.17	0.75
1:AA:15:G:H2'	1:AA:16:A:C8	2.21	0.75
1:AA:675:A:H1'	11:AK:116:HIS:CE1	2.22	0.75
13:AM:90:LEU:HA	13:AM:93:ARG:HB2	1.68	0.75
16:AP:27:LYS:H	16:AP:27:LYS:HD2	1.52	0.75
24:AX:21:G:H2'	24:AX:22:U:C5	2.21	0.75
27:B1:83:GLU:HG3	27:B1:86:SER:H	1.52	0.75
33:B7:25:PRO:HA	33:B7:28:ARG:NH2	2.02	0.75
38:BD:255:LYS:HE3	38:BD:256:GLY:N	2.02	0.75
40:BF:84:VAL:O	40:BF:86:GLY:N	2.19	0.75
50:BT:11:GLU:N	50:BT:11:GLU:CD	2.39	0.75
1:CA:1240:U:H3	7:CG:30:ILE:CG2	1.99	0.75
3:CC:155:GLY:HA3	3:CC:163:ALA:HB1	1.69	0.75
10:CJ:4:ILE:HB	10:CJ:74:ILE:CD1	2.15	0.75
16:CP:19:ILE:HG22	16:CP:36:ILE:HG13	1.68	0.75
35:DA:1644:C:O2'	35:DA:1645:G:H5'	1.87	0.75
38:DD:210:GLY:O	38:DD:211:ARG:HB3	1.87	0.75
41:DG:48:GLU:HG2	41:DG:49:ASP:H	1.52	0.75
42:DH:67:LEU:O	42:DH:71:LEU:HD13	1.87	0.75
56:DZ:150:LEU:HD23	56:DZ:171:ILE:HD11	1.67	0.75
1:AA:1431:C:C2'	1:AA:1432:G:H5'	2.17	0.75
1:AA:1466:C:H2'	1:AA:1467:G:O4'	1.86	0.75
19:AS:6:LYS:HD2	19:AS:6:LYS:H	1.50	0.75
23:AW:24:C:H6	23:AW:24:C:H5'	1.51	0.75
35:BA:1677:A:H2'	35:BA:1678:G:C8	2.22	0.75
35:BA:2195:C:O2'	35:BA:2196:C:H5'	1.87	0.75
35:BA:2300:G:H1	35:BA:2316:C:H42	1.35	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:87:ILE:HD12	45:BO:91:LEU:C	2.06	0.75
56:BZ:149:SER:CB	56:BZ:173:ALA:HA	2.16	0.75
7:CG:100:ALA:O	7:CG:104:LEU:HD23	1.87	0.75
1:CA:600:C:OP1	8:CH:97:VAL:HG12	1.85	0.75
25:CY:179:LYS:O	25:CY:182:GLU:HB2	1.86	0.75
27:D1:78:LYS:HD2	27:D1:78:LYS:N	2.02	0.75
29:D3:52:HIS:CD2	36:DB:83:G:H4'	2.21	0.75
35:DA:2316:C:H1'	41:DG:128:ARG:NH1	2.02	0.75
35:DA:286:C:H2'	35:DA:286:C:O2	1.86	0.75
39:DE:176:ILE:HB	39:DE:181:LEU:HD23	1.67	0.75
41:DG:4:ASP:HA	41:DG:8:LYS:HD3	1.69	0.75
48:DR:24:GLN:HB2	48:DR:44:LEU:HD21	1.66	0.75
54:DX:51:VAL:HG13	54:DX:81:VAL:H	1.52	0.75
1:AA:1184:G:H2'	1:AA:1185:G:C8	2.22	0.75
1:AA:1264:C:H2'	1:AA:1265:G:H8	1.51	0.75
6:AF:53:ALA:HB3	6:AF:86:ARG:HH12	1.49	0.75
16:AP:26:ARG:NH1	16:AP:26:ARG:HB3	2.02	0.75
35:BA:2475:C:H42	35:BA:2529:G:N2	1.84	0.75
35:BA:361:G:H2'	35:BA:362:U:H5''	1.68	0.75
41:BG:43:LEU:HD23	41:BG:44:GLY:H	1.49	0.75
54:BX:40:LYS:HD2	54:BX:41:ASN:N	2.02	0.75
54:BX:47:PHE:O	54:BX:49:VAL:HG23	1.87	0.75
56:BZ:29:TYR:HE2	56:BZ:87:ASP:HB2	1.52	0.75
1:CA:737:A:H2'	1:CA:738:C:H6	1.47	0.75
14:CN:16:PHE:N	14:CN:16:PHE:CD2	2.54	0.75
19:CS:51:VAL:O	19:CS:57:HIS:HA	1.86	0.75
35:DA:2661:G:H2'	35:DA:2662:A:C8	2.22	0.75
39:DE:61:ARG:HG2	39:DE:62:PRO:HD3	1.69	0.75
43:DI:49:ALA:HA	43:DI:52:ARG:HG2	1.68	0.75
5:AE:152:ARG:HB3	8:AH:43:GLY:HA3	1.67	0.74
6:AF:60:PHE:C	6:AF:61:LEU:HD12	2.08	0.74
6:AF:82:ARG:HB2	6:AF:85:VAL:HG23	1.69	0.74
27:B1:20:ARG:NH2	27:B1:41:ARG:HE	1.85	0.74
34:B8:32:LEU:HG	34:B8:34:TRP:CE3	2.22	0.74
35:BA:1418:G:H1	35:BA:1579:A:H5'	1.52	0.74
35:BA:919:G:H4'	36:BB:81:G:O2'	1.86	0.74
42:BH:122:THR:HB	42:BH:134:SER:HB2	1.69	0.74
54:BX:78:LYS:HD3	54:BX:78:LYS:O	1.87	0.74
56:BZ:134:PRO:O	56:BZ:136:PHE:N	2.19	0.74
4:CD:90:GLY:HA3	4:CD:204:ILE:HD11	1.69	0.74
17:CQ:54:GLY:HA3	17:CQ:82:MET:SD	2.27	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:74:ARG:HE	18:CR:81:PHE:HA	1.51	0.74
25:CY:76:LEU:HD21	25:CY:99:LEU:HD21	1.68	0.74
32:D6:11:LEU:HD11	32:D6:51:GLU:HB2	1.67	0.74
35:DA:1484:G:C2'	35:DA:1485:G:H5''	2.16	0.74
35:DA:2475:C:H42	35:DA:2529:G:N2	1.85	0.74
35:DA:832:G:O3'	46:DP:45:LEU:HD11	1.86	0.74
43:DI:133:HIS:HB2	43:DI:134:PRO:HD3	1.66	0.74
35:DA:1245:G:H3'	46:DP:16:ARG:HH22	1.50	0.74
35:DA:587:C:C5	46:DP:33:ARG:HD2	2.22	0.74
45:DO:79:PHE:HA	50:DT:72:VAL:HG22	1.69	0.74
50:DT:62:THR:HG21	50:DT:75:ILE:HG13	1.69	0.74
52:DV:5:VAL:HG21	52:DV:36:PRO:HB2	1.67	0.74
53:DW:13:SER:HB3	53:DW:16:LYS:HD2	1.68	0.74
1:AA:328:C:H4'	1:AA:329:A:H5'	1.69	0.74
1:AA:972:C:H4'	10:AJ:57:LYS:CG	2.17	0.74
2:AB:169:LYS:C	2:AB:169:LYS:HD3	2.07	0.74
2:AB:185:ILE:HG22	2:AB:199:TYR:HB2	1.68	0.74
5:AE:102:ALA:HB2	5:AE:120:THR:OG1	1.87	0.74
5:AE:35:GLY:CA	5:AE:41:VAL:HG12	2.17	0.74
6:AF:33:TYR:CD1	6:AF:75:LEU:HG	2.23	0.74
9:AI:107:ARG:O	9:AI:108:VAL:HG13	1.87	0.74
13:AM:10:PRO:HG2	13:AM:18:ALA:HB1	1.69	0.74
18:AR:74:ARG:HE	18:AR:81:PHE:HA	1.52	0.74
26:B0:21:LEU:HD11	26:B0:41:ARG:HD3	1.68	0.74
35:BA:1550:C:H2'	35:BA:1551:C:H6	1.51	0.74
35:BA:1639:U:C2'	35:BA:1640:C:H5''	2.16	0.74
35:BA:2102:U:H2'	35:BA:2103:C:C6	2.22	0.74
35:BA:797:C:H2'	35:BA:798:G:C8	2.23	0.74
39:BE:102:VAL:HA	39:BE:200:GLU:HA	1.68	0.74
35:BA:2811:G:OP1	39:BE:60:ASN:HB2	1.87	0.74
47:BQ:43:THR:OG1	47:BQ:46:GLN:HG3	1.87	0.74
52:BV:32:THR:HG22	52:BV:33:VAL:N	2.02	0.74
55:BY:31:LEU:HD12	55:BY:33:LYS:N	2.00	0.74
3:CC:150:LYS:HG3	3:CC:169:ALA:HB2	1.69	0.74
5:CE:76:ILE:HG12	5:CE:77:PRO:HD2	1.68	0.74
5:CE:90:VAL:HG23	5:CE:121:LYS:HB3	1.67	0.74
8:CH:119:LEU:HB2	8:CH:123:GLU:HB2	1.68	0.74
1:CA:675:A:H1'	11:CK:116:HIS:CE1	2.22	0.74
13:CM:46:LYS:HG3	13:CM:47:ASP:H	1.52	0.74
35:DA:1019:U:H2'	35:DA:1020:A:C8	2.22	0.74
35:DA:1290:C:H2'	35:DA:1291:C:H6	1.50	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1484:G:C3'	35:DA:1485:G:H5''	2.16	0.74
35:DA:191:A:O2'	35:DA:192:C:H5'	1.87	0.74
35:DA:71:A:H4'	35:DA:72:U:O5'	1.85	0.74
38:DD:255:LYS:HE3	38:DD:256:GLY:N	2.02	0.74
43:DI:115:ALA:H	43:DI:131:LYS:HE2	1.51	0.74
43:DI:115:ALA:N	43:DI:131:LYS:HE2	2.02	0.74
44:DN:43:THR:HB	44:DN:46:VAL:HB	1.67	0.74
44:DN:58:ASP:C	44:DN:60:ILE:H	1.86	0.74
46:DP:45:LEU:HD23	46:DP:46:LYS:H	1.50	0.74
55:DY:15:VAL:HG12	55:DY:16:ALA:H	1.52	0.74
2:AB:71:VAL:HG21	2:AB:93:VAL:HG23	1.68	0.74
6:AF:18:GLN:HA	6:AF:21:LEU:HD23	1.70	0.74
12:AL:119:LYS:HB2	12:AL:120:TYR:CD1	2.22	0.74
16:AP:8:ARG:HG2	16:AP:8:ARG:HH11	1.52	0.74
35:BA:2728:U:O2'	35:BA:2729:G:H5'	1.88	0.74
35:BA:729:G:C5	38:BD:208:LYS:HB2	2.22	0.74
26:B0:77:ARG:HH22	35:BA:857:C:H5'	1.50	0.74
40:BF:110:LEU:HD23	40:BF:110:LEU:O	1.87	0.74
52:BV:3:ALA:HB3	52:BV:14:VAL:HB	1.69	0.74
52:BV:38:LEU:CD2	52:BV:40:LEU:H	2.00	0.74
1:CA:15:G:H2'	1:CA:16:A:C8	2.21	0.74
1:CA:939:G:C5'	7:CG:102:ARG:HH22	2.00	0.74
3:CC:112:SER:HB3	3:CC:115:LEU:CD1	2.17	0.74
11:CK:84:VAL:HG23	11:CK:110:ASP:OD1	1.87	0.74
18:CR:53:ARG:HA	18:CR:56:THR:OG1	1.87	0.74
1:CA:1222:G:H5''	19:CS:78:ARG:NH1	2.01	0.74
23:CW:59:A:H1'	23:CW:61:U:C5	2.22	0.74
35:DA:1528(A):A:H3'	35:DA:1529:G:H5''	1.69	0.74
35:DA:2850:A:H2'	35:DA:2851:A:C8	2.21	0.74
26:D0:77:ARG:HH22	35:DA:857:C:H5'	1.52	0.74
36:DB:66:A:H61	36:DB:108:U:H2'	1.52	0.74
41:DG:144:ILE:CD1	41:DG:145:THR:H	2.00	0.74
44:DN:120:LEU:CD1	44:DN:122:VAL:HG23	2.17	0.74
35:DA:251:A:H5''	46:DP:51:PHE:CZ	2.22	0.74
1:AA:1073:U:H2'	1:AA:1074:G:H8	1.49	0.74
1:AA:148:G:H2'	1:AA:149:A:H8	1.51	0.74
1:AA:920:U:H2'	1:AA:921:U:C6	2.23	0.74
4:AD:28:SER:HB3	4:AD:29:PRO:CD	2.17	0.74
20:AT:38:LYS:O	20:AT:41:ILE:HG12	1.87	0.74
24:AX:21:G:H2'	24:AX:22:U:C6	2.22	0.74
25:AY:5:GLU:O	25:AY:9:GLU:HG3	1.85	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:196:A:H5''	46:BP:46:LYS:HZ1	1.50	0.74
35:BA:2103:C:C3'	35:BA:2104:G:H5''	2.17	0.74
35:BA:832:G:O3'	46:BP:45:LEU:HD11	1.87	0.74
41:BG:173:LEU:HD12	41:BG:178:PHE:CE2	2.23	0.74
42:BH:149:ARG:HA	42:BH:162:ILE:HD11	1.70	0.74
47:BQ:85:LYS:HG3	47:BQ:86:GLY:N	2.02	0.74
56:BZ:103:ARG:HG3	56:BZ:136:PHE:CE1	2.22	0.74
2:CB:185:ILE:HG22	2:CB:199:TYR:HB2	1.68	0.74
2:CB:69:LEU:HD22	2:CB:91:PRO:HB2	1.68	0.74
17:CQ:69:LYS:O	17:CQ:70:ARG:HD2	1.86	0.74
25:CY:115:VAL:O	25:CY:118:VAL:HG23	1.88	0.74
27:D1:94:LEU:CD1	27:D1:95:LEU:H	2.00	0.74
35:DA:1835:G:H5'	35:DA:1836:C:OP2	1.87	0.74
35:DA:2464:C:HO2'	35:DA:2465:C:H6	1.33	0.74
35:DA:404:C:C4'	35:DA:405:U:H5'	2.18	0.74
35:DA:524:U:H2'	35:DA:524:U:O2	1.87	0.74
35:DA:540:C:H2'	35:DA:541:C:C6	2.22	0.74
35:DA:696:G:O2'	35:DA:697:C:H5'	1.87	0.74
40:DF:176:LEU:HD21	40:DF:180:GLY:O	1.87	0.74
44:DN:34:LEU:CD2	44:DN:120:LEU:HB2	2.16	0.74
46:DP:39:LYS:HD2	46:DP:40:SER:H	1.51	0.74
1:AA:1244:C:H2'	1:AA:1245:A:C8	2.22	0.74
1:AA:313:A:H2'	1:AA:314:C:C6	2.22	0.74
1:AA:376:G:OP1	16:AP:6:LEU:HD13	1.86	0.74
20:AT:100:ILE:HD12	20:AT:100:ILE:H	1.52	0.74
27:B1:13:ILE:HG13	27:B1:14:VAL:N	2.00	0.74
35:BA:1403:C:H5''	35:BA:1471:A:C1'	2.17	0.74
35:BA:1484:G:C2'	35:BA:1485:G:H5''	2.16	0.74
35:BA:1876:A:H2'	35:BA:1877:A:C8	2.22	0.74
39:BE:177:PRO:HG2	39:BE:178:GLU:OE1	1.87	0.74
47:BQ:34:LEU:HD12	47:BQ:35:VAL:H	1.52	0.74
48:BR:11:ASN:O	48:BR:12:ARG:HG3	1.87	0.74
49:BS:13:ARG:N	49:BS:13:ARG:HD2	2.03	0.74
49:BS:83:LYS:HA	49:BS:104:GLY:HA2	1.70	0.74
56:BZ:6:LYS:HE3	56:BZ:6:LYS:H	1.52	0.74
1:CA:686:U:H1'	1:CA:687:A:N7	2.01	0.74
6:CF:73:ASN:O	6:CF:76:ALA:HB3	1.87	0.74
1:CA:376:G:OP1	16:CP:6:LEU:HD13	1.87	0.74
26:D0:49:LYS:HB2	26:D0:80:HIS:HB3	1.68	0.74
27:D1:40:ARG:HG2	27:D1:41:ARG:N	2.01	0.74
34:D8:39:LYS:HG2	34:D8:42:ARG:NH1	2.03	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2639:A:H2'	35:DA:2640:G:C5'	2.14	0.74
35:DA:2820:A:C8	39:DE:191:PRO:HB2	2.23	0.74
41:DG:166:ASP:O	41:DG:170:ARG:HB2	1.87	0.74
41:DG:32:PRO:CB	41:DG:172:LEU:HD22	2.17	0.74
43:DI:11:ASN:ND2	43:DI:12:LEU:HD22	2.03	0.74
48:DR:24:GLN:HE22	48:DR:36:THR:HG21	1.50	0.74
49:DS:89:ARG:HA	49:DS:89:ARG:NE	2.03	0.74
52:DV:28:GLU:HB2	52:DV:29:PRO:CD	2.10	0.74
1:AA:179:A:H2'	1:AA:180:U:C6	2.23	0.74
1:AA:137:C:H42	1:AA:226:G:H1	1.36	0.74
4:AD:26:CYS:HA	4:AD:31:CYS:HB2	1.69	0.74
5:AE:57:LYS:O	5:AE:61:TYR:HB2	1.87	0.74
1:AA:939:G:C5'	7:AG:102:ARG:HH22	1.99	0.74
11:AK:84:VAL:HG23	11:AK:110:ASP:OD1	1.87	0.74
17:AQ:3:LYS:HD2	17:AQ:60:ILE:HD11	1.69	0.74
20:AT:18:GLN:O	20:AT:22:ARG:HG3	1.88	0.74
23:AW:10:G:H22	23:AW:27:G:H1'	1.49	0.74
26:B0:27:GLU:N	26:B0:69:PHE:HE1	1.85	0.74
35:BA:286:C:O2	35:BA:286:C:H2'	1.88	0.74
35:BA:342:G:O2'	35:BA:343:C:H5''	1.86	0.74
43:BI:56:LYS:HA	43:BI:59:ALA:HB3	1.68	0.74
46:BP:90:ARG:HD2	46:BP:91:PHE:HD1	1.53	0.74
1:CA:1074:G:H2'	1:CA:1075:C:C6	2.22	0.74
1:CA:1271:G:H5'	1:CA:1314:C:H5''	1.67	0.74
1:CA:1508:G:H2'	1:CA:1509:C:C6	2.21	0.74
6:CF:82:ARG:HB2	6:CF:85:VAL:HG23	1.68	0.74
23:CW:9:G:H2'	23:CW:10:G:N7	2.02	0.74
32:D6:11:LEU:HG	32:D6:26:ASN:HD21	1.51	0.74
35:DA:1428:C:O2'	35:DA:1429:G:H5'	1.87	0.74
43:DI:87:LYS:NZ	43:DI:121:LYS:HG2	2.03	0.74
46:DP:84:ASN:ND2	46:DP:116:GLY:HA3	2.02	0.74
54:DX:8:ILE:H	54:DX:8:ILE:HD12	1.50	0.74
1:AA:1432:G:OP1	50:BT:107:ASP:HB2	1.87	0.74
1:AA:383:A:C2'	1:AA:384:G:H5'	2.18	0.74
2:AB:69:LEU:HD12	2:AB:71:VAL:HG23	1.70	0.74
4:AD:61:LYS:CA	4:AD:203:VAL:HG22	2.18	0.74
6:AF:99:ALA:HB3	18:AR:29:PHE:CE2	2.23	0.74
35:BA:132:G:H5'	35:BA:132:G:H8	1.52	0.74
31:B5:2:ALA:HA	35:BA:2015:A:C1'	2.18	0.74
35:BA:549:G:C2'	35:BA:551:G:H5''	2.17	0.74
45:BO:35:VAL:HG21	45:BO:69:ILE:HG12	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:115:LEU:HA	46:BP:134:ALA:CB	2.17	0.74
48:BR:17:ARG:HH11	48:BR:17:ARG:HG2	1.51	0.74
54:BX:73:ARG:H	54:BX:74:PRO:HD3	1.53	0.74
1:CA:665:A:H2'	1:CA:725:G:H22	1.50	0.74
2:CB:53:ARG:O	2:CB:56:ARG:HB2	1.88	0.74
35:DA:1176:G:H1'	35:DA:1177:A:OP1	1.87	0.74
35:DA:2575:C:H5''	39:DE:144:ARG:HD3	1.70	0.74
35:DA:292:C:H2'	35:DA:293:U:H5'	1.70	0.74
38:DD:229:VAL:HG23	38:DD:230:ASP:H	1.51	0.74
39:DE:197:ILE:HG13	39:DE:199:ARG:HH12	1.53	0.74
40:DF:6:VAL:HG21	40:DF:124:LEU:HA	1.70	0.74
1:AA:976:G:N2	1:AA:1362:C:H2'	2.02	0.74
1:AA:1190:G:H8	3:AC:3:ASN:ND2	1.86	0.74
10:AJ:34:VAL:CG2	10:AJ:74:ILE:HG22	2.16	0.74
10:AJ:84:GLN:O	10:AJ:88:LEU:HB2	1.88	0.74
35:BA:2307:G:H3'	35:BA:2307:G:N3	2.02	0.74
35:BA:259:G:H21	35:BA:621:A:H8	1.31	0.74
35:BA:2691:C:H5'	35:BA:2691:C:H6	1.52	0.74
35:BA:292:C:H2'	35:BA:293:U:H5'	1.70	0.74
38:BD:25:THR:CG2	38:BD:82:ILE:H	2.00	0.74
40:BF:3:GLU:HB2	40:BF:24:LEU:HG	1.67	0.74
45:BO:119:PRO:HB2	50:BT:68:TYR:CE1	2.22	0.74
1:CA:1072:G:H2'	1:CA:1073:U:C6	2.22	0.74
1:CA:376:G:OP1	16:CP:5:ARG:HB2	1.87	0.74
1:CA:806:C:H2'	1:CA:807:A:C8	2.20	0.74
4:CD:17:VAL:HG12	4:CD:18:LYS:H	1.52	0.74
4:AD:197:PRO:HD3	6:CF:16:GLN:HG3	1.69	0.74
8:CH:48:TYR:HA	8:CH:60:ARG:O	1.88	0.74
11:CK:111:ASP:HA	18:CR:84:LYS:HG3	1.68	0.74
13:CM:10:PRO:HG2	13:CM:18:ALA:HB1	1.68	0.74
27:D1:9:GLY:N	27:D1:48:LYS:NZ	2.36	0.74
43:DI:133:HIS:ND1	43:DI:134:PRO:HD2	2.03	0.74
28:D2:23:LYS:CA	54:DX:5:TYR:HE1	2.01	0.74
1:AA:1240:U:H3	7:AG:30:ILE:CG2	2.01	0.74
1:AA:385:C:O2'	1:AA:386:C:H5'	1.88	0.74
1:AA:417:C:O2'	1:AA:418:C:H5'	1.87	0.74
1:AA:735:C:H2'	1:AA:736:C:H6	1.53	0.74
2:AB:96:ARG:CD	2:AB:96:ARG:H	1.89	0.74
4:AD:80:GLU:O	4:AD:84:LYS:HG2	1.88	0.74
1:AA:1222:G:H5''	19:AS:78:ARG:NH1	2.03	0.74
35:BA:1007:C:O2'	44:BN:108:PRO:HA	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2681:C:H5	35:BA:2725:A:H62	1.36	0.74
36:BB:105:A:O2'	56:BZ:30:ASN:HA	1.88	0.74
41:BG:101:ILE:CG1	41:BG:105:LYS:HE3	2.17	0.74
49:BS:92:TYR:CD1	49:BS:93:LYS:N	2.56	0.74
1:CA:1488:G:H2'	1:CA:1489:G:H8	1.51	0.74
1:CA:148:G:H2'	1:CA:149:A:H8	1.53	0.74
1:CA:1512:U:H2'	1:CA:1513:A:H8	1.52	0.74
6:CF:18:GLN:HA	6:CF:21:LEU:HD23	1.68	0.74
6:CF:50:TYR:CE1	18:CR:77:GLY:HA2	2.22	0.74
11:CK:21:ILE:HB	11:CK:84:VAL:HG12	1.68	0.74
25:CY:84:ARG:NH2	25:CY:92:PRO:HD2	2.03	0.74
35:DA:144:C:H2'	35:DA:145:G:C8	2.22	0.74
35:DA:549:G:C2'	35:DA:551:G:H5''	2.18	0.74
35:DA:852:G:O2'	35:DA:853:G:H5'	1.87	0.74
41:DG:43:LEU:H	41:DG:43:LEU:HD13	1.52	0.74
46:DP:115:LEU:HA	46:DP:134:ALA:CB	2.18	0.74
47:DQ:85:LYS:HG3	47:DQ:86:GLY:N	2.01	0.74
49:DS:25:ARG:O	49:DS:39:ILE:HA	1.88	0.74
50:DT:13:ARG:C	50:DT:14:TYR:HD1	1.90	0.74
35:DA:1011:G:OP1	51:DU:75:ASN:HB2	1.87	0.74
51:DU:83:LEU:HB3	51:DU:88:ILE:HG12	1.69	0.74
55:DY:31:LEU:HB2	55:DY:36:ALA:H	1.52	0.74
56:DZ:25:PRO:HG2	56:DZ:85:HIS:HA	1.69	0.74
1:AA:1190:G:P	3:AC:5:ILE:HG23	2.27	0.74
16:AP:18:ARG:HD3	16:AP:35:LYS:HD2	1.68	0.74
18:AR:86:VAL:HG12	18:AR:87:ARG:HH12	1.52	0.74
35:BA:1412:A:H2'	35:BA:1413:G:C8	2.23	0.74
35:BA:2502:G:H5'	35:BA:2503:A:H5''	1.69	0.74
35:BA:272(C):G:H2'	35:BA:272(D):G:H8	1.53	0.74
41:BG:11:TYR:HA	41:BG:15:VAL:HG21	1.70	0.74
47:BQ:35:VAL:HG23	47:BQ:102:VAL:HA	1.68	0.74
47:BQ:8:LYS:HG3	47:BQ:9:TYR:N	2.03	0.74
50:BT:88:ILE:HG22	50:BT:89:VAL:N	2.02	0.74
51:BU:88:ILE:C	51:BU:90:VAL:N	2.40	0.74
53:BW:87:PRO:HA	53:BW:93:ALA:HB2	1.68	0.74
56:BZ:61:LEU:HB2	56:BZ:65:GLN:HB3	1.70	0.74
47:BQ:141:GLN:O	56:BZ:70:LEU:HD22	1.87	0.74
1:CA:1190:G:H8	3:CC:3:ASN:ND2	1.85	0.74
1:CA:1244:C:H2'	1:CA:1245:A:C8	2.23	0.74
1:CA:1317:C:OP1	14:CN:17:LYS:HG2	1.87	0.74
1:CA:383:A:C2'	1:CA:384:G:H5'	2.16	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:69:LEU:HD12	2:CB:71:VAL:HG23	1.68	0.74
18:CR:52:PRO:O	18:CR:56:THR:HG23	1.86	0.74
35:DA:1607:C:H4'	35:DA:1608:A:O5'	1.87	0.74
35:DA:227:A:C2	35:DA:2407:G:H1'	2.23	0.74
35:DA:914:C:H2'	35:DA:915:C:H5'	1.69	0.74
40:DF:110:LEU:HD23	40:DF:110:LEU:O	1.87	0.74
48:DR:103:ARG:HG2	48:DR:103:ARG:HH11	1.51	0.74
52:DV:38:LEU:CD2	52:DV:40:LEU:H	2.00	0.74
52:DV:46:VAL:HG12	52:DV:47:VAL:HG12	1.70	0.74
28:D2:23:LYS:HA	54:DX:5:TYR:HE1	1.52	0.74
1:AA:600:C:O2'	1:AA:601:C:H5'	1.87	0.73
1:AA:673:G:H5''	6:AF:87:ARG:HH11	1.52	0.73
27:B1:78:LYS:N	27:B1:78:LYS:HD2	2.03	0.73
27:B1:86:SER:CA	27:B1:89:GLU:HG3	2.16	0.73
34:B8:32:LEU:O	34:B8:33:ASN:HB3	1.85	0.73
35:BA:1401:G:H2'	35:BA:1402:C:H6	1.53	0.73
35:BA:2661:G:H2'	35:BA:2662:A:C8	2.23	0.73
38:BD:177:LEU:HD12	38:BD:181:GLU:HG3	1.69	0.73
51:BU:70:ARG:HA	51:BU:74:LEU:O	1.88	0.73
55:BY:26:LYS:HG2	55:BY:27:VAL:H	1.52	0.73
1:CA:741:G:H2'	1:CA:742:G:H8	1.53	0.73
8:CH:82:HIS:HD2	8:CH:138:TRP:NE1	1.86	0.73
10:CJ:34:VAL:CG2	10:CJ:74:ILE:HG22	2.16	0.73
15:CO:54:ARG:O	15:CO:58:MET:HG3	1.88	0.73
16:CP:8:ARG:HG2	16:CP:8:ARG:HH11	1.51	0.73
27:D1:38:SER:HB3	35:DA:2080:G:H4'	1.70	0.73
35:DA:1755:A:H2	35:DA:2716:U:H1'	1.51	0.73
35:DA:676:A:H8	35:DA:2069:G:H21	1.35	0.73
36:DB:11:C:H3'	36:DB:12:C:C6	2.23	0.73
39:DE:116:VAL:HG22	39:DE:117:MET:N	2.03	0.73
49:DS:85:VAL:HG23	49:DS:86:ALA:H	1.53	0.73
1:AA:665:A:H2'	1:AA:725:G:H22	1.53	0.73
2:AB:69:LEU:HD22	2:AB:91:PRO:HB2	1.68	0.73
18:AR:37:VAL:HG23	18:AR:38:GLU:N	2.01	0.73
18:AR:53:ARG:HH12	18:AR:59:SER:HA	1.52	0.73
28:B2:29:LYS:O	28:B2:33:MET:N	2.20	0.73
35:BA:1024:G:C3'	35:BA:1025:G:H5''	2.19	0.73
35:BA:2394:C:OP1	46:BP:63:PRO:HD2	1.88	0.73
35:BA:2857:G:N2	35:BA:2859:G:H3'	2.03	0.73
35:BA:588:U:O5'	35:BA:588:U:H6	1.70	0.73
41:BG:169:ALA:O	41:BG:173:LEU:HD23	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BN:14:VAL:HG12	44:BN:15:LEU:N	2.03	0.73
1:CA:1264:C:H2'	1:CA:1265:G:H8	1.52	0.73
1:CA:922:G:N3	1:CA:1398:A:H2	1.86	0.73
1:CA:1442(A):G:C3'	1:CA:1442(B):A:H5''	2.10	0.73
1:CA:303:A:OP1	12:CL:17:LYS:HE3	1.87	0.73
2:CB:49:GLU:O	2:CB:52:GLU:HB3	1.88	0.73
4:CD:132:ARG:HD2	4:CD:132:ARG:O	1.88	0.73
10:CJ:84:GLN:O	10:CJ:88:LEU:HB2	1.86	0.73
17:CQ:88:TYR:O	17:CQ:91:ARG:HB3	1.87	0.73
26:D0:70:GLN:CG	26:D0:71:ASP:H	2.01	0.73
35:DA:690:G:H2'	35:DA:691:C:H6	1.53	0.73
35:DA:1819:A:OP1	38:DD:161:THR:HG21	1.88	0.73
39:DE:177:PRO:HG2	39:DE:178:GLU:OE1	1.89	0.73
40:DF:29:ASN:H	40:DF:112:MET:HE3	1.51	0.73
43:DI:88:ILE:HD11	43:DI:123:LEU:HD12	1.68	0.73
55:DY:75:ILE:HD13	55:DY:76:CYS:N	2.01	0.73
1:AA:551:U:H2'	1:AA:552:U:C6	2.23	0.73
2:AB:29:ALA:O	2:AB:31:TYR:N	2.20	0.73
3:AC:111:LEU:HD21	3:AC:146:ALA:HB2	1.70	0.73
4:AD:30:LYS:C	4:AD:32:ALA:N	2.39	0.73
5:AE:35:GLY:HA3	5:AE:41:VAL:HG12	1.71	0.73
11:AK:17:GLY:HA3	11:AK:80:VAL:HA	1.70	0.73
12:AL:46:LYS:HG2	12:AL:47:LYS:N	2.04	0.73
12:AL:47:LYS:HB3	12:AL:48:PRO:CD	2.18	0.73
14:AN:8:GLU:HB2	14:AN:12:ARG:NH1	2.03	0.73
1:AA:658:G:H1'	15:AO:22:THR:HB	1.69	0.73
35:BA:2291:U:H3	35:BA:2341:G:H1	1.35	0.73
40:BF:103:LYS:HG2	40:BF:106:ARG:NH2	2.02	0.73
40:BF:112:MET:HA	40:BF:115:ALA:HB3	1.70	0.73
43:BI:49:ALA:HA	43:BI:52:ARG:HG2	1.68	0.73
45:BO:13:ASN:O	45:BO:15:GLY:N	2.20	0.73
51:BU:83:LEU:HB3	51:BU:88:ILE:HG12	1.68	0.73
28:B2:29:LYS:NZ	54:BX:9:LEU:HA	2.04	0.73
56:BZ:110:GLY:O	56:BZ:112:ARG:N	2.21	0.73
56:BZ:40:ASP:HB3	56:BZ:43:GLU:HB2	1.70	0.73
1:CA:1443:G:H22	1:CA:1460:A:H1'	1.51	0.73
1:CA:559:A:H4'	1:CA:560:U:C5'	2.18	0.73
2:CB:111:ARG:HG2	2:CB:111:ARG:HH11	1.52	0.73
1:CA:694:A:H5''	11:CK:53:SER:HB3	1.69	0.73
27:D1:13:ILE:HG23	27:D1:14:VAL:H	1.50	0.73
32:D6:15:GLU:OE2	32:D6:41:PRO:HG3	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:D7:34:ARG:HB3	33:D7:42:LEU:HD23	1.70	0.73
34:D8:14:VAL:CG1	34:D8:22:VAL:HG13	2.18	0.73
38:DD:133:LEU:HA	38:DD:136:ILE:HD12	1.68	0.73
38:DD:25:THR:CB	38:DD:82:ILE:H	2.00	0.73
39:DE:9:VAL:HG13	39:DE:25:VAL:O	1.88	0.73
44:DN:137:LYS:HG2	44:DN:138:LEU:H	1.53	0.73
51:DU:70:ARG:HA	51:DU:74:LEU:O	1.89	0.73
55:DY:31:LEU:HD12	55:DY:33:LYS:N	2.03	0.73
1:AA:522:C:N4	1:AA:528:C:H42	1.86	0.73
2:AB:53:ARG:O	2:AB:56:ARG:HB2	1.89	0.73
7:AG:100:ALA:O	7:AG:104:LEU:HD23	1.87	0.73
27:B1:45:ASN:C	27:B1:46:LEU:HD12	2.09	0.73
35:BA:120:U:H1'	35:BA:149:A:C8	2.23	0.73
35:BA:1578:U:H3'	35:BA:1579:A:H5''	1.71	0.73
35:BA:191:A:O2'	35:BA:192:C:H5'	1.87	0.73
36:BB:66:A:H61	36:BB:108:U:H2'	1.52	0.73
41:BG:174:GLU:HA	41:BG:178:PHE:HB2	1.70	0.73
43:BI:11:ASN:ND2	43:BI:12:LEU:HD22	2.03	0.73
43:BI:88:ILE:CG2	43:BI:89:TYR:N	2.50	0.73
46:BP:131:SER:HB2	46:BP:134:ALA:HB3	1.68	0.73
48:BR:95:THR:HA	48:BR:116:LEU:O	1.88	0.73
55:BY:28:LYS:HB2	55:BY:37:VAL:HB	1.70	0.73
56:BZ:125:LEU:HD23	56:BZ:126:VAL:H	1.53	0.73
1:CA:16:A:O2'	1:CA:17:U:H5'	1.88	0.73
1:CA:406:G:H1	1:CA:436:C:H42	1.32	0.73
10:CJ:9:ARG:O	10:CJ:94:VAL:HG13	1.88	0.73
12:CL:86:ARG:HG2	12:CL:87:GLY:H	1.52	0.73
14:CN:8:GLU:HB2	14:CN:12:ARG:NH1	2.03	0.73
35:DA:2300:G:H1	35:DA:2316:C:H42	1.34	0.73
39:DE:167:VAL:HG22	39:DE:168:MET:N	2.03	0.73
42:DH:84:SER:O	42:DH:85:LYS:HB2	1.86	0.73
46:DP:90:ARG:HD2	46:DP:91:PHE:HD1	1.53	0.73
48:DR:18:LEU:HD13	48:DR:19:ALA:N	2.02	0.73
54:DX:47:PHE:O	54:DX:49:VAL:HG23	1.87	0.73
55:DY:26:LYS:HG2	55:DY:27:VAL:H	1.52	0.73
56:DZ:151:HIS:HA	56:DZ:170:THR:HA	1.71	0.73
1:AA:1428:A:H2'	1:AA:1429:C:C6	2.23	0.73
1:AA:428:G:H4'	1:AA:429:U:O5'	1.87	0.73
2:AB:98:LEU:HB2	2:AB:101:MET:HG2	1.71	0.73
4:AD:17:VAL:HG12	4:AD:18:LYS:H	1.52	0.73
6:AF:14:LEU:CD1	6:AF:19:LEU:HB2	2.19	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:44:SER:H	11:AK:47:VAL:HG21	1.52	0.73
14:AN:16:PHE:HD2	14:AN:16:PHE:N	1.86	0.73
23:AW:37:U:H2'	23:AW:38:A:C8	2.23	0.73
25:AY:170:ALA:C	25:AY:172:ALA:H	1.91	0.73
26:B0:36:ILE:HD11	35:BA:2355:C:H4'	1.70	0.73
41:BG:111:LEU:HB2	41:BG:112:PRO:HD3	1.71	0.73
41:BG:170:ARG:HH22	41:BG:182:LYS:CE	2.01	0.73
47:BQ:16:ARG:NH1	47:BQ:16:ARG:HB3	2.03	0.73
53:BW:11:ARG:NH2	53:BW:98:LYS:HB3	2.02	0.73
31:B5:25:LEU:HB2	53:BW:23:LEU:HD11	1.70	0.73
1:CA:1116:C:H2'	1:CA:1117:G:C5'	2.17	0.73
1:CA:892:A:H2'	1:CA:893:C:C6	2.23	0.73
14:CN:16:PHE:N	14:CN:16:PHE:HD2	1.87	0.73
33:D7:34:ARG:HD3	33:D7:42:LEU:HA	1.70	0.73
34:D8:32:LEU:HG	34:D8:34:TRP:CE3	2.22	0.73
34:D8:8:LYS:O	34:D8:12:LYS:HG3	1.88	0.73
26:D0:41:ARG:HH21	35:DA:2387:U:C1'	2.01	0.73
41:DG:130:ASN:O	41:DG:159:VAL:HG23	1.87	0.73
42:DH:92:ILE:HG22	42:DH:93:GLY:N	2.04	0.73
54:DX:76:ARG:O	54:DX:76:ARG:HD3	1.89	0.73
1:AA:410:G:H21	1:AA:432:A:H62	1.35	0.73
3:AC:112:SER:HB3	3:AC:115:LEU:CD1	2.18	0.73
7:AG:137:LYS:O	7:AG:141:VAL:HG23	1.88	0.73
35:BA:1937:A:O2'	35:BA:1938:A:H5'	1.89	0.73
35:BA:97:C:H2'	35:BA:98:G:C8	2.24	0.73
35:BA:992:C:H2'	35:BA:993:G:H8	1.51	0.73
41:BG:60:LEU:CD2	41:BG:63:ILE:HD11	2.12	0.73
42:BH:84:SER:O	42:BH:85:LYS:HB2	1.88	0.73
44:BN:23:LEU:HA	44:BN:26:LEU:HB3	1.70	0.73
28:B2:23:LYS:HA	54:BX:5:TYR:HE1	1.52	0.73
1:CA:403:C:H2'	1:CA:404:U:H6	1.52	0.73
1:CA:971:G:H4'	1:CA:972:C:H5''	1.70	0.73
17:CQ:46:ASP:OD1	17:CQ:49:GLU:HA	1.89	0.73
35:DA:2307:G:N2	35:DA:2308:G:H5'	2.03	0.73
35:DA:2677:G:H2'	35:DA:2678:C:H6	1.52	0.73
35:DA:2704:C:H2'	35:DA:2705:A:C8	2.23	0.73
35:DA:610:G:H22	35:DA:619:G:H1'	1.54	0.73
41:DG:103:LEU:HA	41:DG:106:LEU:HD23	1.70	0.73
52:DV:28:GLU:CB	52:DV:29:PRO:HD3	2.14	0.73
56:DZ:146:ILE:HG22	56:DZ:174:VAL:HG12	1.71	0.73
56:DZ:150:LEU:HD22	56:DZ:150:LEU:H	1.54	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:46:LYS:NZ	12:AL:47:LYS:HB2	2.04	0.73
19:AS:53:ASN:HD21	19:AS:56:GLN:N	1.86	0.73
22:AV:38:U:H2'	22:AV:39:C:C6	2.24	0.73
32:B6:51:GLU:HG2	32:B6:52:VAL:N	2.01	0.73
35:BA:2715:C:H2'	35:BA:2716:U:C6	2.23	0.73
35:BA:2762:G:H2'	35:BA:2763:G:H5''	1.71	0.73
38:BD:229:VAL:HG23	38:BD:230:ASP:H	1.54	0.73
39:BE:181:LEU:N	39:BE:181:LEU:HD22	2.03	0.73
39:BE:61:ARG:HG2	39:BE:62:PRO:HD3	1.69	0.73
45:BO:22:ILE:HB	45:BO:40:VAL:HG12	1.70	0.73
46:BP:83:VAL:CG1	46:BP:112:LEU:HD21	2.18	0.73
50:BT:65:LYS:HZ2	50:BT:65:LYS:HA	1.52	0.73
51:BU:90:VAL:HG22	52:BV:39:LEU:HD12	1.70	0.73
53:BW:32:ALA:O	53:BW:35:ILE:HB	1.88	0.73
54:BX:72:LYS:CG	54:BX:74:PRO:HD3	2.08	0.73
1:CA:1388:C:H2'	1:CA:1389:C:C6	2.24	0.73
1:CA:1522:U:O2'	1:CA:1523:G:H5'	1.89	0.73
1:CA:337:C:H2'	1:CA:338:A:H8	1.54	0.73
4:CD:58:LEU:HD13	4:CD:59:ARG:N	2.02	0.73
9:CI:114:TYR:N	9:CI:114:TYR:CD2	2.50	0.73
12:CL:85:ILE:HD11	12:CL:98:TYR:CB	2.17	0.73
35:DA:154:G:H1	35:DA:172:C:H42	1.36	0.73
35:DA:1550:C:H2'	35:DA:1551:C:H6	1.52	0.73
35:DA:1639:U:O2'	35:DA:1640:C:H5''	1.88	0.73
35:DA:2854:G:H2'	35:DA:2855:C:C6	2.23	0.73
36:DB:91:C:H2'	36:DB:92:C:H6	1.53	0.73
38:DD:209:ALA:O	38:DD:212:SER:HB3	1.88	0.73
39:DE:172:VAL:HG13	39:DE:182:LEU:HD11	1.69	0.73
41:DG:58:GLN:HG3	41:DG:59:GLU:N	2.03	0.73
43:DI:37:VAL:HG13	43:DI:38:LEU:HD12	1.68	0.73
43:DI:56:LYS:HA	43:DI:59:ALA:HB3	1.69	0.73
35:DA:869:G:H1'	47:DQ:8:LYS:NZ	2.04	0.73
55:DY:68:HIS:HB3	55:DY:71:LYS:NZ	2.03	0.73
1:AA:1317:C:OP1	14:AN:17:LYS:HG2	1.88	0.73
8:AH:11:THR:CG2	8:AH:14:ARG:HH12	2.02	0.73
35:BA:389:G:N1	46:BP:71:VAL:HB	2.01	0.73
35:BA:860:U:H5	35:BA:917:A:N7	1.85	0.73
41:BG:33:ARG:HD3	41:BG:162:THR:HG21	1.71	0.73
49:BS:25:ARG:O	49:BS:39:ILE:HA	1.89	0.73
50:BT:13:ARG:C	50:BT:14:TYR:HD1	1.92	0.73
50:BT:38:ASN:ND2	50:BT:40:THR:H	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:48:ILE:O	50:BT:63:VAL:HA	1.88	0.73
55:BY:28:LYS:HZ3	55:BY:30:VAL:HA	1.54	0.73
56:BZ:29:TYR:HE2	56:BZ:87:ASP:CB	2.01	0.73
56:BZ:59:LEU:O	56:BZ:66:SER:HA	1.88	0.73
2:CB:98:LEU:HB2	2:CB:101:MET:HG2	1.69	0.73
6:CF:60:PHE:C	6:CF:61:LEU:HD12	2.09	0.73
19:CS:6:LYS:HD2	19:CS:6:LYS:H	1.53	0.73
27:D1:17:SER:O	27:D1:44:PRO:HD2	1.89	0.73
34:D8:6:THR:CG2	34:D8:63:PRO:HD3	2.18	0.73
35:DA:1449:A:N3	35:DA:1529:G:H1'	2.03	0.73
35:DA:361:G:H2'	35:DA:362:U:H5''	1.70	0.73
39:DE:30:PRO:HD3	39:DE:180:ASN:ND2	2.02	0.73
40:DF:114:VAL:HG23	40:DF:115:ALA:N	2.03	0.73
40:DF:170:LEU:HD12	40:DF:171:PRO:HD2	1.68	0.73
41:DG:137:GLU:HG2	41:DG:138:GLN:H	1.53	0.73
42:DH:73:ALA:O	42:DH:76:VAL:HB	1.88	0.73
44:DN:22:THR:CA	44:DN:61:ARG:HB2	2.19	0.73
50:DT:38:ASN:ND2	50:DT:40:THR:H	1.85	0.73
56:DZ:108:PRO:CB	56:DZ:144:LEU:H	2.01	0.73
1:AA:1489:G:H2'	1:AA:1490:C:C6	2.24	0.73
1:AA:20:U:H2'	1:AA:21:G:O4'	1.87	0.73
1:AA:662:G:H2'	1:AA:663:A:H8	1.51	0.73
12:AL:75:HIS:HD2	12:AL:77:LEU:HG	1.52	0.73
22:AV:39:C:O2'	22:AV:40:C:H5'	1.89	0.73
35:BA:1644:C:O2'	35:BA:1645:G:H5'	1.89	0.73
38:BD:239:ARG:HH21	38:BD:239:ARG:HG3	1.53	0.73
43:BI:54:GLN:HG2	43:BI:57:ARG:HH12	1.53	0.73
45:BO:6:THR:HG22	45:BO:7:TYR:N	2.03	0.73
45:BO:107:ARG:HH11	50:BT:36:GLU:H	1.37	0.73
52:BV:46:VAL:HG12	52:BV:47:VAL:HG12	1.69	0.73
53:BW:51:LEU:HD13	53:BW:52:GLU:N	2.04	0.73
54:BX:5:TYR:O	54:BX:7:VAL:N	2.21	0.73
4:CD:61:LYS:CA	4:CD:203:VAL:HG22	2.18	0.73
1:CA:658:G:C1'	15:CO:22:THR:HB	2.19	0.73
20:CT:43:LEU:HD12	20:CT:52:ALA:HA	1.71	0.73
22:CV:41:C:H2'	22:CV:42:C:C6	2.23	0.73
26:D0:21:LEU:HD11	26:D0:41:ARG:HD3	1.70	0.73
27:D1:83:GLU:HG2	27:D1:86:SER:HB2	1.71	0.73
35:DA:2092:U:H5	35:DA:2226:C:OP2	1.72	0.73
35:DA:271(U):G:H2'	35:DA:271(V):G:C8	2.24	0.73
35:DA:705:A:H1'	38:DD:9:TYR:CE1	2.24	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:95:ARG:NH1	41:DG:95:ARG:HA	2.04	0.73
1:AA:498:U:O2	1:AA:498:U:H2'	1.89	0.73
3:AC:157:ILE:HB	3:AC:164:ARG:NH1	2.04	0.73
4:AD:142:PRO:HA	4:AD:185:PHE:HD2	1.54	0.73
8:AH:119:LEU:HB2	8:AH:123:GLU:HB2	1.71	0.73
23:AW:42:C:H2'	23:AW:43:G:H8	1.54	0.73
27:B1:14:VAL:O	27:B1:46:LEU:HD23	1.89	0.73
35:BA:1258:C:H2'	35:BA:1259:G:C8	2.23	0.73
35:BA:1614:A:C2	53:BW:87:PRO:HB3	2.24	0.73
35:BA:185:U:H2'	35:BA:186:G:H8	1.53	0.73
35:BA:2007:C:H2'	35:BA:2008:C:H6	1.52	0.73
35:BA:571:A:H5'	35:BA:2030:A:H62	1.54	0.73
38:BD:45:ASN:CG	38:BD:46:GLN:H	1.89	0.73
38:BD:25:THR:HG22	38:BD:82:ILE:O	1.89	0.73
39:BE:77:ILE:HG22	39:BE:78:LEU:N	2.04	0.73
42:BH:105:LEU:HD22	42:BH:105:LEU:N	2.03	0.73
42:BH:67:LEU:O	42:BH:71:LEU:HD13	1.87	0.73
43:BI:54:GLN:HG2	43:BI:57:ARG:NH1	2.03	0.73
46:BP:84:ASN:ND2	46:BP:116:GLY:HA3	2.04	0.73
23:CW:2:G:N2	23:CW:3:C:H1'	2.03	0.73
28:D2:44:LEU:C	28:D2:46:GLN:H	1.92	0.73
29:D3:56:VAL:HG12	29:D3:57:GLU:N	2.03	0.73
31:D5:42:PRO:HB2	35:DA:2815:C:O2'	1.89	0.73
35:DA:2487:G:H2'	35:DA:2488:A:H8	1.54	0.73
35:DA:2801(A):A:C4'	35:DA:2802:G:H5'	2.19	0.73
35:DA:342:G:O2'	35:DA:343:C:H5''	1.87	0.73
35:DA:2773:C:H5''	39:DE:164:ARG:O	1.89	0.73
41:DG:137:GLU:HG2	41:DG:138:GLN:N	2.03	0.73
50:DT:22:PHE:N	50:DT:22:PHE:CD2	2.51	0.73
52:DV:70:ILE:CB	52:DV:90:PRO:HB2	2.17	0.73
1:AA:1442(B):A:OP1	1:AA:1442(B):A:H4'	1.88	0.72
1:AA:154:C:H2'	1:AA:155:C:C6	2.24	0.72
2:AB:133:LYS:O	2:AB:137:ARG:HB2	1.88	0.72
4:AD:120:LEU:N	4:AD:120:LEU:HD12	2.01	0.72
8:AH:48:TYR:HA	8:AH:60:ARG:O	1.88	0.72
11:AK:21:ILE:N	11:AK:21:ILE:HD12	2.04	0.72
35:BA:208:C:H2'	35:BA:209:C:C6	2.23	0.72
26:B0:41:ARG:HH21	35:BA:2387:U:C1'	2.01	0.72
35:BA:676:A:H8	35:BA:2069:G:H21	1.34	0.72
39:BE:82:ARG:HG3	39:BE:83:ASP:H	1.54	0.72
45:BO:69:ILE:HD13	45:BO:77:ILE:HG23	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BV:4:ILE:HD12	52:BV:40:LEU:HD21	1.71	0.72
55:BY:74:PRO:O	55:BY:75:ILE:HB	1.87	0.72
1:CA:1303:C:O2	1:CA:1303:C:H2'	1.88	0.72
1:CA:1435:G:H2'	1:CA:1436:U:C6	2.24	0.72
6:CF:76:ALA:HB1	6:CF:80:ARG:NH2	2.04	0.72
16:CP:18:ARG:HD3	16:CP:35:LYS:HD2	1.71	0.72
28:D2:14:ARG:CG	28:D2:15:LYS:N	2.50	0.72
28:D2:40:SER:CB	28:D2:41:ILE:HD12	2.19	0.72
35:DA:272(C):G:H2'	35:DA:272(D):G:H8	1.54	0.72
35:DA:648:G:H2'	35:DA:649:G:H8	1.53	0.72
41:DG:76:SER:HB3	41:DG:83:ARG:CB	2.19	0.72
42:DH:153:LYS:HB2	42:DH:154:PRO:HD2	1.71	0.72
45:DO:16:ALA:HB1	45:DO:43:VAL:HG13	1.71	0.72
46:DP:83:VAL:HG12	46:DP:112:LEU:HD21	1.71	0.72
54:DX:54:VAL:C	54:DX:55:ASN:HD22	1.93	0.72
1:AA:1522:U:O2'	1:AA:1523:G:H5'	1.89	0.72
1:AA:985:C:H2'	1:AA:986:A:C8	2.24	0.72
4:AD:31:CYS:O	4:AD:31:CYS:SG	2.46	0.72
20:AT:60:GLU:HG3	20:AT:81:LYS:HE3	1.71	0.72
7:AG:148:ASN:ND2	23:AW:41:C:H4'	2.04	0.72
35:BA:144:C:H2'	35:BA:145:G:C8	2.24	0.72
35:BA:1509(B):A:H2'	35:BA:1510:G:C8	2.24	0.72
35:BA:1607:C:H4'	35:BA:1608:A:O5'	1.88	0.72
35:BA:1689:A:H62	35:BA:1698:A:H2	1.34	0.72
35:BA:1908:C:H2'	35:BA:1909:C:H6	1.53	0.72
35:BA:2558:C:H2'	35:BA:2559:C:C6	2.25	0.72
35:BA:296:C:O2'	35:BA:297:C:H5'	1.89	0.72
38:BD:135:PHE:HD1	38:BD:135:PHE:H	1.35	0.72
44:BN:132:ALA:O	44:BN:133:GLN:HB2	1.87	0.72
46:BP:47:ASP:HB3	46:BP:48:PRO:O	1.88	0.72
50:BT:102:ILE:HA	50:BT:110:ILE:HD11	1.71	0.72
56:BZ:127:LYS:H	56:BZ:164:ALA:HB3	1.49	0.72
1:CA:186:C:H2'	1:CA:187:C:C6	2.24	0.72
8:CH:32:LYS:O	8:CH:35:ILE:HG12	1.89	0.72
9:CI:28:VAL:HG13	9:CI:63:ILE:O	1.89	0.72
20:CT:18:GLN:O	20:CT:22:ARG:HG3	1.88	0.72
35:DA:2562:U:H2'	35:DA:2563:U:H5'	1.70	0.72
35:DA:2570:G:H2'	35:DA:2571:C:C6	2.25	0.72
35:DA:2821:A:H2'	35:DA:2822:G:C8	2.23	0.72
35:DA:541:C:H2'	35:DA:542:C:C6	2.23	0.72
39:DE:98:PRO:HG3	39:DE:175:VAL:HG12	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:70:THR:O	42:DH:72:ILE:N	2.22	0.72
47:DQ:141:GLN:HE21	56:DZ:72:ARG:HA	1.54	0.72
47:DQ:27:VAL:HG22	56:DZ:81:ARG:HH22	1.53	0.72
49:DS:83:LYS:HA	49:DS:104:GLY:HA2	1.69	0.72
56:DZ:44:PHE:CE2	56:DZ:86:VAL:HG21	2.23	0.72
1:AA:553:A:H2'	1:AA:554:C:C6	2.25	0.72
3:AC:149:ALA:HA	3:AC:201:TYR:O	1.90	0.72
8:AH:120:THR:HG23	8:AH:123:GLU:OE1	1.88	0.72
9:AI:9:ARG:HG3	9:AI:14:VAL:HG22	1.69	0.72
25:AY:117:ALA:O	25:AY:120:GLN:HB3	1.88	0.72
28:B2:14:ARG:HE	28:B2:14:ARG:N	1.86	0.72
34:B8:52:LYS:H	34:B8:53:PRO:CD	1.99	0.72
35:BA:1503:U:H2'	35:BA:1504:C:C6	2.24	0.72
35:BA:329:G:OP2	55:BY:71:LYS:HD3	1.89	0.72
35:BA:481:G:H1'	35:BA:506:G:N2	2.04	0.72
38:BD:183:ARG:HG2	38:BD:183:ARG:HH11	1.53	0.72
40:BF:8:GLN:HB2	40:BF:126:VAL:HA	1.71	0.72
1:CA:706:A:C5	1:CA:707:C:H5	2.08	0.72
1:CA:854:G:H3'	1:CA:871:U:O4	1.89	0.72
2:CB:133:LYS:O	2:CB:137:ARG:HB2	1.88	0.72
9:CI:107:ARG:O	9:CI:108:VAL:HG13	1.88	0.72
11:CK:44:SER:N	11:CK:47:VAL:HG21	2.04	0.72
16:CP:27:LYS:H	16:CP:27:LYS:HD2	1.54	0.72
28:D2:37:PHE:CE2	28:D2:40:SER:HA	2.24	0.72
35:DA:1876:A:H2'	35:DA:1877:A:C8	2.24	0.72
35:DA:2000:G:HO2'	35:DA:2689:U:H5	1.36	0.72
35:DA:2521:C:H42	35:DA:2544:G:H1	1.37	0.72
35:DA:992:C:H2'	35:DA:993:G:H8	1.53	0.72
38:DD:14:ARG:HH11	38:DD:14:ARG:CB	2.01	0.72
39:DE:59:VAL:HG11	39:DE:63:LEU:HG	1.70	0.72
1:AA:186:C:H2'	1:AA:187:C:C6	2.24	0.72
1:AA:266:G:H22	1:AA:270:A:H62	1.38	0.72
4:AD:153:ARG:HB3	4:AD:153:ARG:NH1	2.05	0.72
9:AI:50:LEU:O	9:AI:53:VAL:HG22	1.89	0.72
10:AJ:82:ILE:O	10:AJ:86:MET:HB3	1.89	0.72
23:AW:28:U:H3	23:AW:44:A:H2	1.37	0.72
27:B1:83:GLU:CG	27:B1:86:SER:H	2.02	0.72
35:BA:1341:U:P	35:BA:1397:U:H3	2.11	0.72
35:BA:227:A:C2	35:BA:2407:G:H1'	2.25	0.72
35:BA:71:A:H4'	35:BA:72:U:O5'	1.89	0.72
42:BH:105:LEU:HD21	42:BH:113:VAL:HB	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BQ:8:LYS:CG	47:BQ:9:TYR:H	2.01	0.72
54:BX:55:ASN:HB2	54:BX:77:LYS:HD2	1.72	0.72
1:CA:20:U:H2'	1:CA:21:G:O4'	1.87	0.72
1:CA:985:C:H2'	1:CA:986:A:C8	2.25	0.72
4:CD:26:CYS:HA	4:CD:31:CYS:HB2	1.71	0.72
5:CE:51:VAL:O	5:CE:55:VAL:HG23	1.89	0.72
27:D1:83:GLU:CG	27:D1:86:SER:HB2	2.19	0.72
35:DA:1170:G:H1	35:DA:1179:C:H42	1.37	0.72
35:DA:1660:C:H5'	35:DA:2712(A):A:H61	1.54	0.72
35:DA:342:G:H2'	35:DA:343:C:C5'	2.20	0.72
35:DA:467:G:H2'	35:DA:468:G:H8	1.52	0.72
39:DE:82:ARG:HG3	39:DE:83:ASP:H	1.53	0.72
41:DG:144:ILE:HG13	41:DG:145:THR:N	2.04	0.72
47:DQ:51:ARG:O	47:DQ:54:MET:HB3	1.89	0.72
56:DZ:166:SER:CB	56:DZ:168:GLU:N	2.40	0.72
1:AA:1303:C:H2'	1:AA:1303:C:O2	1.88	0.72
1:AA:1346:A:N1	1:AA:1374:A:H5''	2.05	0.72
1:AA:1404:C:H2'	1:AA:1405:G:C8	2.24	0.72
1:AA:971:G:H4'	1:AA:972:C:H5''	1.70	0.72
4:AD:90:GLY:HA3	4:AD:204:ILE:HD11	1.72	0.72
5:AE:145:LYS:O	5:AE:148:VAL:HB	1.89	0.72
28:B2:17:SER:O	28:B2:21:LEU:N	2.21	0.72
33:B7:9:ARG:NH1	35:BA:1309:G:H3'	2.05	0.72
35:BA:1270:C:H5''	35:BA:1271:G:H5'	1.71	0.72
35:BA:1290:C:H2'	35:BA:1291:C:C6	2.24	0.72
35:BA:1342:A:H5'	54:BX:55:ASN:OD1	1.89	0.72
35:BA:2074:U:H2'	35:BA:2075:U:C6	2.25	0.72
35:BA:2720:U:H5'	35:BA:2721:A:OP2	1.89	0.72
28:B2:55:ARG:NH1	35:BA:72:U:H5'	2.04	0.72
38:BD:129:ASN:O	38:BD:193:VAL:HG12	1.89	0.72
40:BF:10:PRO:HD2	40:BF:13:SER:O	1.88	0.72
40:BF:6:VAL:HG21	40:BF:124:LEU:HA	1.72	0.72
41:BG:27:ASN:HD21	41:BG:29:TRP:CB	1.99	0.72
50:BT:109:GLU:O	50:BT:113:LYS:HG3	1.88	0.72
50:BT:52:ILE:O	50:BT:52:ILE:HG13	1.88	0.72
54:BX:64:LYS:O	54:BX:65:ARG:HB2	1.88	0.72
1:CA:1118:C:H6	1:CA:1118:C:H5'	1.54	0.72
1:CA:17:U:H2'	1:CA:18:C:C6	2.24	0.72
2:CB:178:ARG:NH2	8:CH:74:PRO:HG3	2.03	0.72
3:CC:9:GLY:O	3:CC:12:LEU:HB2	1.89	0.72
4:CD:28:SER:HB3	4:CD:29:PRO:CD	2.17	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:90:VAL:CG2	5:CE:121:LYS:HB3	2.20	0.72
8:CH:120:THR:OG1	8:CH:123:GLU:HG3	1.88	0.72
25:CY:36:ALA:HA	25:CY:39:LEU:HG	1.71	0.72
26:D0:27:GLU:HB2	26:D0:69:PHE:CD1	2.25	0.72
27:D1:13:ILE:CG1	27:D1:14:VAL:N	2.50	0.72
27:D1:14:VAL:O	27:D1:46:LEU:HD23	1.89	0.72
31:D5:56:LYS:O	31:D5:57:VAL:HG13	1.89	0.72
35:DA:1345:C:H2'	35:DA:1346:G:H8	1.53	0.72
35:DA:1503:U:H2'	35:DA:1504:C:C6	2.24	0.72
35:DA:2102:U:H2'	35:DA:2103:C:C6	2.23	0.72
35:DA:2852:G:H2'	35:DA:2853:C:C6	2.24	0.72
35:DA:329:G:OP2	55:DY:71:LYS:HD3	1.90	0.72
39:DE:128:SER:O	39:DE:129:HIS:HB2	1.90	0.72
43:DI:102:SER:CB	43:DI:109:ILE:HG12	2.19	0.72
45:DO:87:ILE:HD12	45:DO:91:LEU:C	2.09	0.72
55:DY:37:VAL:HG23	55:DY:38:ILE:N	2.04	0.72
56:DZ:131:ARG:O	56:DZ:133:ILE:HD12	1.88	0.72
2:AB:72:GLY:HA3	2:AB:165:VAL:HG13	1.70	0.72
3:AC:107:GLN:CD	3:AC:107:GLN:H	1.93	0.72
35:BA:1528(A):A:H3'	35:BA:1529:G:H5''	1.70	0.72
35:BA:2189:U:C3'	35:BA:2190:G:H5''	2.19	0.72
35:BA:285:C:H2'	35:BA:286:C:H5''	1.72	0.72
36:BB:11:C:H3'	36:BB:12:C:C6	2.24	0.72
39:BE:77:ILE:CG2	39:BE:78:LEU:N	2.52	0.72
45:BO:114:ILE:H	45:BO:114:ILE:CD1	1.97	0.72
46:BP:122:PRO:HB3	46:BP:141:ALA:HB1	1.72	0.72
47:BQ:33:GLY:HA2	47:BQ:105:GLU:HA	1.70	0.72
51:BU:18:LEU:HD23	51:BU:18:LEU:O	1.89	0.72
54:BX:54:VAL:C	54:BX:55:ASN:HD22	1.92	0.72
1:CA:1080:A:H5'	5:CE:14:ARG:NH2	2.03	0.72
8:CH:37:ARG:O	8:CH:41:ARG:HB3	1.90	0.72
18:CR:47:THR:OG1	18:CR:49:LYS:HG2	1.90	0.72
35:DA:465:G:H2'	35:DA:466:A:C8	2.24	0.72
39:DE:39:PRO:HA	39:DE:43:GLY:HA2	1.69	0.72
50:DT:48:ILE:O	50:DT:63:VAL:HA	1.89	0.72
54:DX:64:LYS:O	54:DX:65:ARG:HB2	1.88	0.72
55:DY:87:LYS:HG3	55:DY:89:PHE:N	2.04	0.72
1:AA:471:G:H2'	1:AA:472:A:H8	1.53	0.72
1:AA:688:G:H2'	1:AA:689:C:H6	1.54	0.72
9:AI:2:GLU:O	9:AI:3:GLN:HG3	1.88	0.72
9:AI:95:LYS:HZ3	9:AI:96:LEU:HB2	1.55	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:88:TYR:O	17:AQ:91:ARG:HB3	1.89	0.72
25:AY:48:ALA:O	25:AY:50:VAL:HG13	1.90	0.72
35:BA:1034:G:N2	35:BA:1122:G:H1'	2.05	0.72
35:BA:585:G:H2'	35:BA:1251:C:H42	1.54	0.72
35:BA:635:C:H2'	35:BA:636:G:C8	2.23	0.72
38:BD:183:ARG:HG3	38:BD:269:PHE:O	1.89	0.72
40:BF:65:TRP:CH2	40:BF:75:HIS:HD2	2.08	0.72
56:BZ:28:MET:HE3	56:BZ:37:VAL:HG21	1.70	0.72
1:CA:1258:G:H2'	1:CA:1259:C:C6	2.25	0.72
1:CA:1445:C:H2'	1:CA:1446:U:H6	1.53	0.72
1:CA:452:A:H4'	16:CP:72:ARG:NH2	2.05	0.72
18:CR:37:VAL:HG23	18:CR:38:GLU:N	2.02	0.72
25:CY:18:LEU:CG	25:CY:19:GLU:N	2.53	0.72
27:D1:58:ILE:HG23	27:D1:59:THR:N	2.03	0.72
28:D2:40:SER:HB2	28:D2:41:ILE:HD12	1.72	0.72
32:D6:51:GLU:HG2	32:D6:52:VAL:N	2.01	0.72
35:DA:185:U:H2'	35:DA:186:G:H8	1.54	0.72
35:DA:610:G:N2	35:DA:619:G:H1'	2.04	0.72
35:DA:635:C:H2'	35:DA:636:G:C8	2.25	0.72
40:DF:132:VAL:HG22	40:DF:133:ASN:N	2.03	0.72
34:D8:30:ARG:NH2	46:DP:62:LEU:HD23	2.04	0.72
50:DT:53:ARG:HG2	50:DT:53:ARG:HH11	1.55	0.72
54:DX:40:LYS:HD2	54:DX:41:ASN:N	2.04	0.72
55:DY:76:CYS:SG	55:DY:77:PRO:HD3	2.28	0.72
1:AA:1118:C:H5'	1:AA:1118:C:H6	1.54	0.72
1:AA:1053:G:O6	1:AA:1200:C:H5''	1.88	0.72
1:AA:444:C:H2'	1:AA:445:G:H8	1.55	0.72
1:AA:522:C:H41	12:AL:53:ARG:NH2	1.85	0.72
1:AA:694:A:O2'	23:AW:39:A:H1'	1.90	0.72
10:AJ:32:ALA:HB1	10:AJ:75:ILE:HD11	1.72	0.72
13:AM:91:ARG:HB3	13:AM:96:LEU:O	1.89	0.72
33:B7:45:ALA:O	33:B7:46:VAL:HB	1.88	0.72
35:BA:1188:U:O2'	35:BA:1189:A:H5'	1.88	0.72
35:BA:2575:C:H5''	39:BE:144:ARG:HD3	1.71	0.72
35:BA:2808:U:H2'	35:BA:2809:A:H5'	1.72	0.72
40:BF:18:ARG:HG2	40:BF:19:GLU:H	1.53	0.72
40:BF:69:HIS:O	40:BF:70:THR:HG23	1.89	0.72
42:BH:103:LEU:HD23	42:BH:115:VAL:HB	1.71	0.72
43:BI:102:SER:CB	43:BI:109:ILE:HG12	2.19	0.72
45:BO:121:VAL:C	45:BO:122:LEU:HD12	2.09	0.72
54:BX:16:LYS:HA	54:BX:16:LYS:HE3	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:45:VAL:CA	55:BY:62:GLU:HG2	2.12	0.72
1:CA:692:U:H2'	1:CA:694:A:OP2	1.89	0.72
1:CA:878:G:H5'	8:CH:89:PRO:CG	2.19	0.72
1:CA:977:A:H2'	1:CA:978:A:H5'	1.72	0.72
1:CA:989:C:H42	1:CA:1217:C:N4	1.88	0.72
2:CB:165:VAL:CG2	2:CB:166:ASP:H	1.99	0.72
12:CL:46:LYS:NZ	12:CL:47:LYS:HB2	2.05	0.72
35:DA:1364:G:H1'	35:DA:1368:G:H22	1.55	0.72
31:D5:2:ALA:HB2	35:DA:2014:A:HO2'	1.55	0.72
35:DA:2052:G:H2'	35:DA:2053:G:H8	1.55	0.72
35:DA:2729:G:H1'	39:DE:187:ALA:CB	2.17	0.72
35:DA:709:U:H2'	35:DA:710:G:H8	1.54	0.72
38:DD:224:ALA:HB2	38:DD:233:HIS:HB3	1.72	0.72
39:DE:77:ILE:HG22	39:DE:78:LEU:N	2.05	0.72
41:DG:73:ALA:HB3	41:DG:87:PRO:HG2	1.71	0.72
47:DQ:76:LYS:H	47:DQ:88:GLY:CA	2.03	0.72
1:AA:375:U:H2'	1:AA:376:G:H8	1.53	0.72
4:AD:3:ARG:O	4:AD:5:ILE:HG13	1.89	0.72
7:AG:71:PRO:HG3	7:AG:103:TRP:CZ3	2.24	0.72
8:AH:6:ILE:HG22	8:AH:10:LEU:HD11	1.72	0.72
9:AI:118:LYS:HB3	9:AI:118:LYS:HZ3	1.53	0.72
19:AS:15:LEU:HD22	19:AS:15:LEU:H	1.54	0.72
27:B1:90:ILE:O	27:B1:93:GLU:HG2	1.89	0.72
28:B2:53:LEU:O	28:B2:54:LYS:HB3	1.89	0.72
35:BA:1131:G:N3	35:BA:1132:A:N7	2.38	0.72
35:BA:1170:G:H1	35:BA:1179:C:H42	1.37	0.72
35:BA:1578:U:C3'	35:BA:1579:A:H5''	2.19	0.72
38:BD:48:ARG:HH11	38:BD:48:ARG:HG3	1.53	0.72
40:BF:176:LEU:HD21	40:BF:180:GLY:O	1.90	0.72
1:CA:542:G:H2'	1:CA:543:C:C6	2.25	0.72
1:CA:728:A:H2'	1:CA:729:A:C8	2.24	0.72
4:CD:33:MET:HE2	4:CD:37:PRO:HA	1.72	0.72
12:CL:47:LYS:HB3	12:CL:48:PRO:CD	2.19	0.72
6:CF:99:ALA:HB3	18:CR:29:PHE:CE2	2.25	0.72
27:D1:38:SER:CB	35:DA:2080:G:H4'	2.20	0.72
29:D3:31:LEU:HD22	29:D3:32:GLN:H	1.55	0.72
35:DA:2307:G:N3	35:DA:2307:G:H3'	2.04	0.72
35:DA:2739:U:O2'	35:DA:2740:A:H5'	1.89	0.72
39:DE:116:VAL:O	39:DE:117:MET:HB3	1.87	0.72
41:DG:52:ILE:O	41:DG:54:GLU:N	2.23	0.72
45:DO:35:VAL:HG21	45:DO:69:ILE:HG12	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:26:LEU:O	49:DS:26:LEU:HD23	1.89	0.72
31:B5:40:LYS:HZ3	31:B5:45:VAL:HA	1.54	0.72
32:B6:36:LEU:HD13	32:B6:50:ARG:NH1	2.05	0.72
35:BA:1449:A:N3	35:BA:1529:G:H1'	2.05	0.72
38:BD:27:THR:CG2	38:BD:28:GLU:H	1.91	0.72
39:BE:154:LYS:HE3	39:BE:154:LYS:HA	1.72	0.72
41:BG:91:ARG:HG2	41:BG:92:VAL:H	1.54	0.72
42:BH:89:ILE:HD13	42:BH:90:LYS:H	1.55	0.72
45:BO:61:VAL:HG22	45:BO:62:VAL:O	1.89	0.72
45:BO:2:ILE:HD11	45:BO:82:ASN:CB	2.19	0.72
46:BP:16:ARG:CD	46:BP:18:ARG:H	2.02	0.72
50:BT:83:ILE:HG13	50:BT:84:GLN:HG2	1.72	0.72
44:BN:41:ASP:C	51:BU:64:ARG:HH11	1.94	0.72
55:BY:2:ARG:N	55:BY:4:LYS:HE2	2.05	0.72
55:BY:2:ARG:N	55:BY:4:LYS:HG2	2.05	0.72
1:CA:1062:U:H2'	1:CA:1063:C:C6	2.24	0.72
1:CA:154:C:H2'	1:CA:155:C:C6	2.25	0.72
7:CG:65:ALA:HB1	7:CG:127:ALA:HB3	1.71	0.72
35:DA:1024:G:C3'	35:DA:1025:G:H5''	2.20	0.72
35:DA:1446:C:H42	35:DA:1465:G:H1	1.38	0.72
35:DA:919:G:H4'	36:DB:81:G:O2'	1.90	0.72
38:DD:267:SER:C	38:DD:269:PHE:N	2.42	0.72
43:DI:77:LEU:HB2	43:DI:140:LEU:CD1	2.17	0.72
44:DN:132:ALA:O	44:DN:133:GLN:HB2	1.88	0.72
46:DP:131:SER:HB2	46:DP:134:ALA:HB3	1.71	0.72
50:DT:109:GLU:HB3	50:DT:113:LYS:CE	2.18	0.72
50:DT:50:ILE:HA	50:DT:99:LEU:HD11	1.72	0.72
52:DV:19:LYS:NZ	52:DV:20:LEU:H	1.87	0.72
54:DX:49:VAL:HG12	54:DX:50:LYS:H	1.54	0.72
55:DY:101:LYS:HG2	55:DY:102:CYS:N	2.05	0.72
55:DY:79:CYS:SG	55:DY:80:GLY:N	2.62	0.72
1:AA:1074:G:H2'	1:AA:1075:C:C6	2.25	0.71
1:AA:114:U:H2'	1:AA:115:G:C8	2.25	0.71
1:AA:1431:C:H2'	1:AA:1432:G:H5'	1.72	0.71
8:AH:37:ARG:O	8:AH:41:ARG:HB3	1.89	0.71
19:AS:51:VAL:O	19:AS:57:HIS:HA	1.89	0.71
29:B3:31:LEU:HD22	29:B3:32:GLN:H	1.55	0.71
35:BA:1786:A:C5	35:BA:1938:A:N7	2.57	0.71
35:BA:226:G:H5'	35:BA:257:A:H4'	1.72	0.71
35:BA:2639:A:H2'	35:BA:2640:G:C5'	2.16	0.71
35:BA:2801(A):A:C4'	35:BA:2802:G:H5'	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2817:G:H21	35:BA:2836:U:H1'	1.52	0.71
35:BA:648:G:H2'	35:BA:649:G:H8	1.55	0.71
38:BD:172:TYR:CD1	38:BD:186:HIS:HA	2.23	0.71
40:BF:164:ARG:HG3	40:BF:175:THR:OG1	1.89	0.71
42:BH:101:ARG:HB2	42:BH:117:PRO:HG3	1.72	0.71
44:BN:107:LEU:HB2	44:BN:108:PRO:HD2	1.71	0.71
48:BR:62:ALA:O	48:BR:66:VAL:HG23	1.90	0.71
52:BV:61:VAL:HB	52:BV:99:ILE:N	2.04	0.71
56:BZ:17:ALA:C	56:BZ:20:ARG:HB3	2.11	0.71
56:BZ:95:PRO:HA	56:BZ:129:SER:HA	1.72	0.71
1:CA:597:G:H2'	1:CA:598:U:H5'	1.72	0.71
3:CC:107:GLN:CD	3:CC:107:GLN:H	1.94	0.71
3:CC:73:PRO:O	3:CC:76:VAL:HG22	1.90	0.71
4:CD:3:ARG:O	4:CD:5:ILE:HG13	1.90	0.71
5:CE:35:GLY:CA	5:CE:41:VAL:HG12	2.20	0.71
5:CE:57:LYS:O	5:CE:61:TYR:HB2	1.89	0.71
27:D1:8:SER:HA	35:DA:1365:A:OP2	1.90	0.71
35:DA:1578:U:H3'	35:DA:1579:A:H5''	1.70	0.71
35:DA:1300:U:O2	35:DA:1626:G:H2'	1.89	0.71
35:DA:2195:C:O2'	35:DA:2196:C:H5'	1.90	0.71
35:DA:2291:U:H3	35:DA:2341:G:H1	1.35	0.71
35:DA:445:C:H5''	51:DU:3:ARG:HB2	1.72	0.71
35:DA:549:G:H2'	35:DA:551:G:C5'	2.19	0.71
35:DA:920:G:H2'	35:DA:921:G:H8	1.54	0.71
38:DD:118:VAL:HG22	38:DD:119:ALA:N	2.05	0.71
38:DD:44:ASN:CB	38:DD:49:ILE:HA	2.18	0.71
41:DG:125:PHE:HD1	41:DG:125:PHE:H	1.37	0.71
41:DG:38:VAL:N	41:DG:158:ALA:HB3	2.05	0.71
42:DH:43:VAL:HG23	42:DH:43:VAL:O	1.89	0.71
43:DI:98:ALA:CB	43:DI:109:ILE:HB	2.19	0.71
43:DI:54:GLN:HG2	43:DI:57:ARG:NH1	2.04	0.71
44:DN:26:LEU:CG	44:DN:30:ILE:HD11	2.19	0.71
46:DP:129:ALA:C	46:DP:130:PHE:HD2	1.93	0.71
1:AA:989:C:H42	1:AA:1217:C:N4	1.89	0.71
5:AE:147:ASP:HA	5:AE:150:ARG:HB3	1.72	0.71
13:AM:46:LYS:HG3	13:AM:47:ASP:H	1.53	0.71
27:B1:13:ILE:HB	27:B1:63:ALA:HB2	1.71	0.71
35:BA:2704:C:H2'	35:BA:2705:A:C8	2.24	0.71
38:BD:267:SER:O	38:BD:269:PHE:N	2.23	0.71
39:BE:197:ILE:HG13	39:BE:199:ARG:HH12	1.55	0.71
35:BA:675:A:H4'	40:BF:67:GLN:OE1	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:104:ARG:HB3	45:BO:104:ARG:NH1	2.06	0.71
51:BU:68:ALA:O	51:BU:71:GLN:HB3	1.88	0.71
55:BY:27:VAL:HG12	55:BY:29:GLU:H	1.55	0.71
56:BZ:118:GLN:O	56:BZ:172:ALA:HA	1.89	0.71
1:CA:1184:G:H2'	1:CA:1185:G:C8	2.24	0.71
1:CA:1241:G:H2'	1:CA:1242:C:C6	2.25	0.71
1:CA:1483:A:H2'	1:CA:1484:C:O4'	1.90	0.71
1:CA:565:U:H3'	1:CA:566:G:H2'	1.71	0.71
1:CA:738:C:H5''	6:CF:69:GLU:HB2	1.71	0.71
7:CG:71:PRO:HG3	7:CG:103:TRP:CZ3	2.26	0.71
12:CL:47:LYS:CG	12:CL:48:PRO:HD3	2.20	0.71
22:CV:30:A:C4	22:CV:31:U:H5	2.08	0.71
25:CY:28:LEU:O	25:CY:30:THR:HG23	1.90	0.71
31:D5:17:ASP:O	31:D5:20:ARG:HB2	1.90	0.71
35:DA:582:G:H2'	35:DA:583:G:C8	2.24	0.71
38:DD:247:ALA:HA	38:DD:254:THR:HG22	1.72	0.71
41:DG:41:GLN:HE22	41:DG:153:ARG:HB3	1.55	0.71
42:DH:101:ARG:HB2	42:DH:117:PRO:HG3	1.72	0.71
45:DO:113:LYS:O	45:DO:117:LEU:HG	1.90	0.71
50:DT:60:THR:HG22	50:DT:77:PRO:HA	1.73	0.71
52:DV:29:PRO:HD2	52:DV:32:THR:HG1	1.55	0.71
53:DW:32:ALA:O	53:DW:35:ILE:HB	1.90	0.71
55:DY:2:ARG:N	55:DY:4:LYS:HG2	2.04	0.71
55:DY:88:LYS:HD2	55:DY:88:LYS:N	2.04	0.71
1:AA:1220:G:H2'	1:AA:1221:G:H8	1.55	0.71
3:AC:155:GLY:HA3	3:AC:163:ALA:HB1	1.70	0.71
8:AH:82:HIS:HD2	8:AH:138:TRP:NE1	1.87	0.71
17:AQ:82:MET:O	17:AQ:85:VAL:HB	1.89	0.71
34:B8:13:ARG:HB3	46:BP:63:PRO:CA	2.15	0.71
35:BA:108:U:H2'	35:BA:109:G:H8	1.56	0.71
35:BA:1314:C:H5'	35:BA:1314:C:H6	1.56	0.71
35:BA:549:G:H2'	35:BA:551:G:C5'	2.19	0.71
35:BA:621:A:H2'	35:BA:622:G:H5'	1.70	0.71
38:BD:79:VAL:CG1	38:BD:113:VAL:HA	2.19	0.71
42:BH:83:TYR:HB3	42:BH:135:GLY:O	1.91	0.71
45:BO:68:GLU:HB3	45:BO:78:ARG:HD3	1.73	0.71
47:BQ:55:VAL:HG12	47:BQ:64:ILE:CD1	2.20	0.71
50:BT:25:GLY:HA2	50:BT:92:GLY:CA	2.20	0.71
54:BX:18:TYR:HA	54:BX:21:PHE:CD1	2.25	0.71
56:BZ:24:LEU:HG	56:BZ:25:PRO:HD2	1.72	0.71
2:CB:233:SER:HB2	2:CB:234:PRO:HD2	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:91:ARG:HB3	13:CM:96:LEU:O	1.90	0.71
17:CQ:3:LYS:HD2	17:CQ:60:ILE:HD11	1.71	0.71
20:CT:100:ILE:H	20:CT:100:ILE:HD12	1.53	0.71
32:D6:36:LEU:HD13	32:D6:50:ARG:NH1	2.05	0.71
35:DA:1221(A):C:O2'	35:DA:1222:C:H5'	1.90	0.71
35:DA:2853:C:H2'	35:DA:2854:G:H8	1.55	0.71
36:DB:7:G:H3'	36:DB:8:U:C5'	2.21	0.71
42:DH:105:LEU:HD22	42:DH:105:LEU:N	2.04	0.71
43:DI:57:ARG:NH1	43:DI:57:ARG:HB3	2.05	0.71
47:DQ:33:GLY:HA2	47:DQ:105:GLU:HA	1.72	0.71
48:DR:95:THR:HA	48:DR:116:LEU:O	1.90	0.71
48:DR:61:HIS:O	48:DR:65:LEU:HB2	1.90	0.71
50:DT:88:ILE:HG22	50:DT:89:VAL:N	2.05	0.71
1:AA:903:G:H2'	1:AA:904:C:H6	1.54	0.71
2:AB:49:GLU:O	2:AB:52:GLU:HB3	1.90	0.71
5:AE:13:ILE:HA	5:AE:29:GLY:O	1.90	0.71
11:AK:21:ILE:HD13	11:AK:84:VAL:HG12	1.72	0.71
12:AL:85:ILE:HD11	12:AL:98:TYR:CB	2.20	0.71
25:AY:76:LEU:CD1	25:AY:99:LEU:HG	2.20	0.71
27:B1:87:PRO:C	27:B1:89:GLU:H	1.92	0.71
29:B3:59:VAL:HG12	29:B3:60:GLU:N	2.03	0.71
35:BA:1788:C:H2'	35:BA:1789:A:H8	1.55	0.71
35:BA:2591:C:OP2	38:BD:239:ARG:HB2	1.90	0.71
35:BA:536:A:H2'	35:BA:537:C:C6	2.25	0.71
35:BA:797:C:H2'	35:BA:798:G:H8	1.53	0.71
39:BE:203:LYS:O	39:BE:203:LYS:HD2	1.90	0.71
41:BG:135:LEU:HG	41:BG:136:ARG:H	1.54	0.71
42:BH:16:SER:CB	42:BH:27:LYS:HB2	2.20	0.71
35:BA:2749:A:H4'	42:BH:62:LYS:HB3	1.73	0.71
44:BN:137:LYS:HG2	44:BN:138:LEU:H	1.56	0.71
47:BQ:25:ASP:HA	56:BZ:78:LYS:HZ1	1.53	0.71
48:BR:24:GLN:HE22	48:BR:36:THR:HG21	1.54	0.71
55:BY:88:LYS:HD2	55:BY:88:LYS:N	2.05	0.71
56:BZ:104:PHE:CD1	56:BZ:139:VAL:HG21	2.25	0.71
1:CA:410:G:H21	1:CA:432:A:H62	1.36	0.71
1:CA:543:C:H2'	1:CA:544:G:C8	2.26	0.71
5:CE:12:LEU:HD13	5:CE:31:LEU:HB3	1.72	0.71
8:CH:120:THR:HG23	8:CH:123:GLU:OE1	1.90	0.71
9:CI:28:VAL:CG1	9:CI:64:THR:HA	2.21	0.71
28:D2:41:ILE:H	28:D2:41:ILE:CD1	1.97	0.71
35:DA:2189:U:C3'	35:DA:2190:G:H5''	2.19	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2201:C:O2'	35:DA:2202:C:H5'	1.90	0.71
35:DA:2749:A:H4'	42:DH:62:LYS:HB3	1.72	0.71
35:DA:2772:C:H2'	35:DA:2773:C:C6	2.25	0.71
40:DF:41:LEU:HD11	40:DF:184:TYR:HE1	1.56	0.71
35:DA:660:G:H5'	40:DF:99:TYR:CE2	2.25	0.71
41:DG:137:GLU:CG	41:DG:138:GLN:H	2.01	0.71
42:DH:122:THR:HB	42:DH:134:SER:HB2	1.71	0.71
45:DO:36:GLY:N	45:DO:62:VAL:HB	2.05	0.71
48:DR:11:ASN:O	48:DR:12:ARG:HG3	1.90	0.71
50:DT:52:ILE:HG13	50:DT:52:ILE:O	1.90	0.71
51:DU:21:ALA:HB2	51:DU:35:ALA:HB1	1.71	0.71
51:DU:93:LYS:HD3	51:DU:93:LYS:H	1.54	0.71
35:DA:1614:A:N6	53:DW:93:ALA:HB2	2.06	0.71
2:AB:105:PHE:O	2:AB:108:ILE:HG22	1.90	0.71
19:AS:43:GLU:C	19:AS:45:VAL:H	1.94	0.71
26:B0:48:GLY:HA3	26:B0:80:HIS:ND1	2.05	0.71
32:B6:11:LEU:HG	32:B6:26:ASN:HD21	1.55	0.71
34:B8:52:LYS:N	34:B8:53:PRO:CD	2.54	0.71
35:BA:2186:G:C2'	35:BA:2187:G:H5''	2.20	0.71
35:BA:2677:G:H2'	35:BA:2678:C:H6	1.55	0.71
35:BA:2859:G:H2'	35:BA:2860:A:C8	2.26	0.71
35:BA:342:G:H2'	35:BA:343:C:C5'	2.19	0.71
35:BA:559:G:N2	51:BU:49:HIS:CD2	2.59	0.71
39:BE:9:VAL:HG13	39:BE:25:VAL:O	1.91	0.71
41:BG:28:VAL:HG12	41:BG:28:VAL:O	1.90	0.71
41:BG:43:LEU:HD21	41:BG:88:ILE:HG22	1.70	0.71
41:BG:46:ALA:HB2	41:BG:88:ILE:HB	1.71	0.71
42:BH:92:ILE:HG22	42:BH:93:GLY:N	2.06	0.71
44:BN:9:VAL:HG11	44:BN:39:ARG:NH2	2.04	0.71
45:BO:2:ILE:HD12	45:BO:6:THR:HG21	1.72	0.71
49:BS:15:ARG:HB3	49:BS:18:ILE:CD1	2.17	0.71
56:BZ:150:LEU:N	56:BZ:150:LEU:HD22	2.05	0.71
1:CA:477:A:O2'	1:CA:479:C:H5'	1.90	0.71
5:CE:72:GLN:HE22	5:CE:77:PRO:HD3	1.55	0.71
9:CI:2:GLU:O	9:CI:3:GLN:HG3	1.90	0.71
10:CJ:30:SER:O	10:CJ:81:THR:HG23	1.91	0.71
12:CL:46:LYS:HG2	12:CL:47:LYS:N	2.04	0.71
35:DA:1412:A:H2'	35:DA:1413:G:C8	2.25	0.71
35:DA:285:C:H2'	35:DA:286:C:H5''	1.72	0.71
35:DA:481:G:H1'	35:DA:506:G:N2	2.06	0.71
35:DA:621:A:H2'	35:DA:622:G:H5'	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:5:VAL:CG1	41:DG:6:ALA:H	1.91	0.71
49:DS:85:VAL:HG23	49:DS:86:ALA:N	2.05	0.71
53:DW:87:PRO:HA	53:DW:93:ALA:HB2	1.73	0.71
55:DY:28:LYS:HB2	55:DY:37:VAL:HB	1.72	0.71
1:AA:1241:G:H2'	1:AA:1242:C:C6	2.26	0.71
1:AA:559:A:H4'	1:AA:560:U:C5'	2.20	0.71
4:AD:132:ARG:HD2	4:AD:132:ARG:O	1.90	0.71
4:AD:33:MET:HE2	4:AD:37:PRO:HA	1.72	0.71
1:AA:738:C:H5''	6:AF:69:GLU:HB2	1.73	0.71
9:AI:83:ARG:O	9:AI:86:VAL:HG12	1.89	0.71
18:AR:78:LEU:O	18:AR:79:LEU:HG	1.90	0.71
35:BA:1258:C:H2'	35:BA:1259:G:H8	1.54	0.71
35:BA:1751:C:O2'	35:BA:1752:C:H5'	1.91	0.71
35:BA:2069:G:O2'	35:BA:2070:G:H5'	1.91	0.71
35:BA:2558:C:H2'	35:BA:2559:C:H6	1.54	0.71
35:BA:2773:C:H5''	39:BE:164:ARG:O	1.89	0.71
39:BE:183:LEU:HD21	50:BT:11:GLU:HB3	1.71	0.71
44:BN:39:ARG:HD3	44:BN:39:ARG:O	1.89	0.71
46:BP:70:GLN:HG3	46:BP:71:VAL:H	1.56	0.71
49:BS:38:GLN:HG2	49:BS:39:ILE:N	2.04	0.71
55:BY:28:LYS:HA	55:BY:39:VAL:H	1.53	0.71
55:BY:88:LYS:HZ3	55:BY:93:GLY:N	1.84	0.71
4:CD:142:PRO:HA	4:CD:185:PHE:HD2	1.55	0.71
26:D0:72:ARG:HD3	26:D0:75:LEU:HD13	1.73	0.71
35:DA:1279:G:H4'	48:DR:31:HIS:CD2	2.24	0.71
35:DA:1860:G:H1	35:DA:1882:C:H42	1.38	0.71
40:DF:112:MET:HA	40:DF:115:ALA:HB3	1.71	0.71
42:DH:83:TYR:HB3	42:DH:135:GLY:O	1.89	0.71
48:DR:62:ALA:O	48:DR:66:VAL:HG23	1.91	0.71
49:DS:15:ARG:HB3	49:DS:18:ILE:CD1	2.14	0.71
50:DT:109:GLU:O	50:DT:113:LYS:HG3	1.89	0.71
1:AA:6:G:H4'	1:AA:298:A:H4'	1.72	0.71
1:AA:542:G:H2'	1:AA:543:C:C6	2.26	0.71
1:AA:706:A:C5	1:AA:707:C:H5	2.09	0.71
3:AC:127:ARG:HG2	3:AC:127:ARG:HH11	1.54	0.71
9:AI:114:TYR:N	9:AI:114:TYR:CD2	2.51	0.71
1:AA:972:C:O3'	10:AJ:57:LYS:HG2	1.90	0.71
10:AJ:42:THR:HG23	10:AJ:68:HIS:HA	1.73	0.71
34:B8:32:LEU:HB3	34:B8:35:GLN:H	1.56	0.71
35:BA:323:G:HO2'	35:BA:1205:U:H3	1.38	0.71
35:BA:792:G:H5''	35:BA:793:A:H5'	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:132:ASN:OD1	41:BG:158:ALA:HA	1.90	0.71
44:BN:120:LEU:CD1	44:BN:122:VAL:HG23	2.21	0.71
45:BO:107:ARG:NH1	50:BT:36:GLU:H	1.89	0.71
46:BP:129:ALA:C	46:BP:130:PHE:HD2	1.93	0.71
46:BP:32:THR:O	46:BP:33:ARG:HB2	1.90	0.71
46:BP:39:LYS:HD3	46:BP:40:SER:H	1.54	0.71
49:BS:85:VAL:HG23	49:BS:86:ALA:N	2.06	0.71
1:CA:1350:A:OP2	9:CI:118:LYS:HD2	1.90	0.71
1:CA:1375:A:H2'	1:CA:1376:U:C6	2.25	0.71
1:CA:15:G:H4'	5:CE:24:ARG:NH2	2.05	0.71
4:CD:59:ARG:O	4:CD:62:GLN:HB2	1.91	0.71
4:CD:5:ILE:HG22	4:CD:6:GLY:N	2.03	0.71
7:CG:151:TYR:O	7:CG:154:TYR:HB2	1.90	0.71
13:CM:9:ILE:HG22	13:CM:11:ARG:HG3	1.72	0.71
33:D7:45:ALA:O	33:D7:46:VAL:HB	1.90	0.71
35:DA:1418:G:H1	35:DA:1579:A:H5'	1.55	0.71
35:DA:285:C:C3'	35:DA:286:C:H5''	2.21	0.71
38:DD:177:LEU:HD12	38:DD:181:GLU:HG3	1.73	0.71
40:DF:10:PRO:HD2	40:DF:13:SER:O	1.90	0.71
41:DG:135:LEU:HD23	41:DG:155:MET:HE1	1.73	0.71
41:DG:144:ILE:HG13	41:DG:145:THR:H	1.54	0.71
43:DI:54:GLN:HG2	43:DI:57:ARG:HH12	1.55	0.71
50:DT:102:ILE:HA	50:DT:110:ILE:HD11	1.72	0.71
50:DT:25:GLY:HA2	50:DT:92:GLY:CA	2.20	0.71
50:DT:83:ILE:HG13	50:DT:84:GLN:N	2.06	0.71
1:AA:955:U:H1'	1:AA:1227:A:N6	2.05	0.71
1:AA:1258:G:H2'	1:AA:1259:C:C6	2.26	0.71
1:AA:1280:A:H5''	10:AJ:40:LEU:HD13	1.73	0.71
1:AA:597:G:H2'	1:AA:598:U:H5'	1.72	0.71
2:AB:178:ARG:HH21	8:AH:74:PRO:HG3	1.56	0.71
2:AB:204:ASN:HD21	2:AB:207:ALA:H	1.39	0.71
2:AB:54:THR:HG22	2:AB:58:ILE:HD11	1.73	0.71
4:AD:128:VAL:HG12	4:AD:129:ASN:N	2.06	0.71
13:AM:23:TYR:CD1	13:AM:67:GLU:HA	2.26	0.71
27:B1:20:ARG:HH12	27:B1:41:ARG:CZ	2.04	0.71
35:BA:1336:A:H2'	35:BA:1337:G:C8	2.25	0.71
35:BA:1364:G:H1'	35:BA:1368:G:H22	1.52	0.71
35:BA:673:C:H5'	40:BF:54:ARG:HH12	1.55	0.71
35:BA:814:C:O2'	35:BA:815:C:H5'	1.90	0.71
45:BO:113:LYS:O	45:BO:117:LEU:HG	1.89	0.71
50:BT:62:THR:HG21	50:BT:75:ILE:HG13	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:BX:49:VAL:HG12	54:BX:50:LYS:H	1.55	0.71
1:CA:266:G:H22	1:CA:270:A:H62	1.39	0.71
2:CB:105:PHE:O	2:CB:108:ILE:HG22	1.91	0.71
1:CA:1349:A:P	9:CI:118:LYS:HZ2	2.12	0.71
19:CS:43:GLU:C	19:CS:45:VAL:H	1.94	0.71
26:D0:48:GLY:HA3	26:D0:80:HIS:ND1	2.05	0.71
35:DA:1034:G:N2	35:DA:1122:G:H1'	2.05	0.71
35:DA:120:U:H1'	35:DA:149:A:C8	2.26	0.71
35:DA:1449:A:C2	35:DA:1529:G:H1'	2.26	0.71
35:DA:2317:C:O2'	35:DA:2318:G:H5'	1.91	0.71
35:DA:2752:C:H5	35:DA:2753:A:H62	1.37	0.71
39:DE:203:LYS:HD2	39:DE:203:LYS:O	1.89	0.71
42:DH:126:PRO:O	42:DH:127:GLU:HG2	1.91	0.71
1:AA:930:C:O2'	1:AA:931:C:H5'	1.91	0.71
11:AK:44:SER:N	11:AK:47:VAL:HG21	2.05	0.71
12:AL:47:LYS:CG	12:AL:48:PRO:HD3	2.20	0.71
26:B0:49:LYS:HB2	26:B0:80:HIS:HB3	1.71	0.71
35:BA:1887:C:H3'	35:BA:1888:G:H5''	1.73	0.71
35:BA:2290:G:H5'	35:BA:2290:G:H8	1.55	0.71
35:BA:285:C:C3'	35:BA:286:C:H5''	2.21	0.71
35:BA:78:A:H2'	35:BA:79:G:H8	1.55	0.71
35:BA:986:C:O2'	35:BA:987:G:H5'	1.91	0.71
36:BB:7:G:H4'	49:BS:29:PHE:CE2	2.25	0.71
38:BD:14:ARG:CB	38:BD:14:ARG:HH11	2.01	0.71
38:BD:210:GLY:O	38:BD:212:SER:N	2.24	0.71
41:BG:178:PHE:HB3	41:BG:180:PHE:CE1	2.24	0.71
44:BN:62:VAL:O	44:BN:63:THR:HG22	1.90	0.71
45:BO:16:ALA:HB1	45:BO:43:VAL:HG13	1.72	0.71
1:CA:1239:A:H62	1:CA:1299:A:H62	1.36	0.71
1:CA:1280:A:H5''	10:CJ:40:LEU:HD13	1.72	0.71
10:CJ:56:HIS:O	10:CJ:58:ASP:N	2.24	0.71
10:CJ:42:THR:HG23	10:CJ:68:HIS:HA	1.72	0.71
25:CY:106:LEU:HD12	25:CY:110:ARG:HG2	1.72	0.71
35:DA:1031:G:H22	35:DA:1124:C:H1'	1.56	0.71
35:DA:2598:A:C5'	38:DD:236:GLY:H	1.94	0.71
39:DE:77:ILE:CG2	39:DE:78:LEU:N	2.53	0.71
45:DO:104:ARG:NH1	45:DO:104:ARG:HB3	2.06	0.71
54:DX:58:HIS:O	54:DX:59:VAL:HG13	1.89	0.71
54:DX:29:TRP:HZ3	54:DX:76:ARG:HG2	1.54	0.71
55:DY:27:VAL:HG12	55:DY:29:GLU:H	1.56	0.71
1:AA:778:G:O2'	1:AA:779:C:H5'	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:150:LYS:HG3	3:AC:169:ALA:HB2	1.72	0.71
3:AC:148:GLY:CA	3:AC:203:PHE:HB3	2.20	0.71
4:AD:119:GLN:HG3	4:AD:123:HIS:HD2	1.54	0.71
13:AM:52:GLU:O	13:AM:56:LEU:HB2	1.91	0.71
13:AM:9:ILE:HG22	13:AM:11:ARG:HG3	1.73	0.71
27:B1:73:LEU:HA	27:B1:76:ARG:HH12	1.56	0.71
35:BA:2853:C:H2'	35:BA:2854:G:H8	1.56	0.71
35:BA:1803:A:O2'	38:BD:259:THR:HG21	1.91	0.71
39:BE:78:LEU:CD2	39:BE:78:LEU:H	2.03	0.71
41:BG:125:PHE:HB2	41:BG:166:ASP:OD2	1.91	0.71
49:BS:28:VAL:HG12	49:BS:29:PHE:N	2.03	0.71
1:CA:417:C:O2'	1:CA:418:C:H5'	1.90	0.71
1:CA:986:A:H1'	19:CS:54:GLY:O	1.91	0.71
2:CB:101:MET:HB2	2:CB:102:LEU:HD12	1.73	0.71
4:CD:59:ARG:HH22	4:CD:66:ARG:NH2	1.89	0.71
4:CD:80:GLU:O	4:CD:84:LYS:HG2	1.91	0.71
12:CL:102:ARG:HG2	12:CL:102:ARG:HH11	1.56	0.71
20:CT:60:GLU:HG3	20:CT:81:LYS:HE3	1.72	0.71
34:D8:52:LYS:HE3	34:D8:52:LYS:HA	1.73	0.71
35:DA:1301:A:H4'	35:DA:1302:A:OP1	1.90	0.71
35:DA:2701:C:H3'	35:DA:2702:U:C5'	2.20	0.71
35:DA:2859:G:H2'	35:DA:2860:A:C8	2.26	0.71
39:DE:101:ARG:HD3	39:DE:169:ASN:ND2	2.05	0.71
41:DG:125:PHE:CD1	41:DG:125:PHE:N	2.57	0.71
41:DG:73:ALA:N	41:DG:87:PRO:HD2	2.06	0.71
45:DO:13:ASN:O	45:DO:15:GLY:N	2.23	0.71
45:DO:107:ARG:HH11	50:DT:36:GLU:H	1.37	0.71
56:DZ:48:PHE:CE2	56:DZ:71:VAL:HG11	2.24	0.71
2:AB:84:GLU:CB	2:AB:219:VAL:HG21	2.20	0.70
3:AC:173:VAL:HG12	3:AC:173:VAL:O	1.91	0.70
11:AK:69:ALA:HA	11:AK:72:ALA:HB3	1.72	0.70
17:AQ:46:ASP:OD1	17:AQ:49:GLU:HA	1.91	0.70
32:B6:32:ASN:CG	32:B6:33:LYS:H	1.94	0.70
34:B8:14:VAL:CG1	34:B8:22:VAL:HG13	2.20	0.70
35:BA:1279:G:H4'	48:BR:31:HIS:CD2	2.26	0.70
35:BA:1345:C:H2'	35:BA:1346:G:H8	1.55	0.70
35:BA:2801(A):A:H4'	35:BA:2802:G:C2'	2.21	0.70
35:BA:807:U:H2'	35:BA:808:G:H8	1.56	0.70
41:BG:172:LEU:N	41:BG:175:LEU:HD12	2.05	0.70
41:BG:47:LYS:HE3	41:BG:81:LYS:CB	2.19	0.70
50:BT:50:ILE:HA	50:BT:99:LEU:HD11	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:68:HIS:HB3	55:BY:71:LYS:NZ	2.06	0.70
56:BZ:58:VAL:HA	56:BZ:67:LEU:O	1.91	0.70
1:CA:1220:G:H2'	1:CA:1221:G:H8	1.55	0.70
4:CD:65:ARG:HD2	4:CD:70:ILE:O	1.91	0.70
6:CF:45:LEU:HD12	6:CF:46:ARG:N	2.06	0.70
16:CP:82:GLN:HE21	16:CP:82:GLN:H	1.39	0.70
27:D1:26:ARG:HA	27:D1:34:THR:HA	1.71	0.70
35:DA:1336:A:H2'	35:DA:1337:G:C8	2.25	0.70
35:DA:1340:U:C6	35:DA:1603:A:O4'	2.44	0.70
35:DA:1342:A:H5'	54:DX:55:ASN:OD1	1.91	0.70
35:DA:2125:G:N2	35:DA:2173:A:H62	1.84	0.70
40:DF:164:ARG:HG3	40:DF:175:THR:OG1	1.90	0.70
42:DH:16:SER:CB	42:DH:27:LYS:HB2	2.21	0.70
56:DZ:33:LEU:HD11	56:DZ:35:ARG:HG2	1.72	0.70
1:AA:1320:C:N4	19:AS:36:ARG:HG3	2.07	0.70
1:AA:1388:C:H2'	1:AA:1389:C:C6	2.26	0.70
1:AA:741:G:H2'	1:AA:742:G:H8	1.56	0.70
1:AA:748:C:H1'	1:AA:749:C:OP2	1.91	0.70
2:AB:101:MET:HA	2:AB:108:ILE:HG13	1.74	0.70
2:AB:69:LEU:CD1	2:AB:71:VAL:HG23	2.20	0.70
7:AG:151:TYR:O	7:AG:154:TYR:HB2	1.90	0.70
12:AL:38:THR:HG22	12:AL:57:LYS:O	1.90	0.70
13:AM:91:ARG:HB3	13:AM:98:VAL:HG22	1.73	0.70
35:BA:1428:C:O2'	35:BA:1429:G:H5'	1.90	0.70
35:BA:1639:U:O2'	35:BA:1640:C:H5''	1.91	0.70
35:BA:2075:U:H2'	35:BA:2238:G:N2	2.06	0.70
39:BE:100:GLU:O	39:BE:172:VAL:HG23	1.90	0.70
39:BE:52:LEU:HD12	39:BE:53:PRO:HD2	1.73	0.70
47:BQ:140:ALA:HB2	56:BZ:99:TYR:CG	2.25	0.70
48:BR:4:LEU:O	48:BR:4:LEU:HD13	1.91	0.70
51:BU:21:ALA:HB2	51:BU:35:ALA:HB1	1.72	0.70
52:BV:28:GLU:CB	52:BV:29:PRO:HD3	2.13	0.70
1:CA:1307:U:H2'	1:CA:1308:U:C6	2.25	0.70
1:CA:1401:G:C2'	1:CA:1402:C:H5'	2.21	0.70
3:CC:173:VAL:O	3:CC:173:VAL:HG12	1.90	0.70
19:CS:15:LEU:H	19:CS:15:LEU:HD22	1.56	0.70
26:D0:23:VAL:HG13	26:D0:37:LEU:O	1.90	0.70
28:D2:49:LYS:O	28:D2:52:ASP:N	2.24	0.70
35:DA:1131:G:N3	35:DA:1132:A:N7	2.39	0.70
35:DA:690:G:H2'	35:DA:691:C:C6	2.26	0.70
35:DA:6:A:O2'	44:DN:130:HIS:HB2	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:172:TYR:CD1	38:DD:186:HIS:HA	2.25	0.70
38:DD:236:GLY:O	38:DD:237:GLU:HG2	1.91	0.70
40:DF:63:LYS:HZ1	40:DF:67:GLN:HB3	1.55	0.70
42:DH:149:ARG:HG3	42:DH:162:ILE:HD11	1.72	0.70
56:DZ:76:LEU:CD2	56:DZ:76:LEU:N	2.54	0.70
1:AA:1057:G:O2'	1:AA:1058:G:H5'	1.91	0.70
1:AA:728:A:H2'	1:AA:729:A:C8	2.25	0.70
1:AA:892:A:H2'	1:AA:893:C:H6	1.55	0.70
1:AA:694:A:H5''	11:AK:53:SER:HB3	1.72	0.70
6:AF:91:VAL:HG11	18:AR:72:ARG:HH12	1.55	0.70
27:B1:87:PRO:O	27:B1:90:ILE:HG12	1.91	0.70
31:B5:56:LYS:O	31:B5:57:VAL:HG13	1.90	0.70
35:BA:6:A:O2'	44:BN:130:HIS:HB2	1.90	0.70
35:BA:740:U:H2'	35:BA:741:G:H8	1.53	0.70
35:BA:803:U:O2'	35:BA:804:A:H5'	1.91	0.70
50:BT:27:THR:HA	50:BT:88:ILE:H	1.53	0.70
56:BZ:162:GLU:O	56:BZ:164:ALA:N	2.24	0.70
56:BZ:17:ALA:HA	56:BZ:20:ARG:HD2	1.71	0.70
1:CA:1443:G:N2	1:CA:1460:A:H1'	2.07	0.70
4:CD:33:MET:CE	4:CD:37:PRO:HA	2.21	0.70
27:D1:16:ASN:ND2	27:D1:16:ASN:H	1.90	0.70
27:D1:67:ILE:N	27:D1:68:PRO:CD	2.55	0.70
33:D7:25:PRO:HG2	33:D7:26:GLY:H	1.56	0.70
35:DA:1290:C:H2'	35:DA:1291:C:C6	2.26	0.70
35:DA:1401:G:H2'	35:DA:1402:C:C6	2.26	0.70
35:DA:1844:C:O2'	35:DA:1845:G:H5'	1.91	0.70
35:DA:2681:C:H5	35:DA:2725:A:H62	1.37	0.70
35:DA:2807:G:H3'	35:DA:2808:U:H5''	1.73	0.70
35:DA:364:C:C2'	35:DA:365:C:H5''	2.20	0.70
36:DB:11:C:H3'	36:DB:12:C:H6	1.56	0.70
41:DG:51:ARG:HE	41:DG:51:ARG:HA	1.56	0.70
41:DG:67:LYS:HD2	41:DG:67:LYS:H	1.56	0.70
44:DN:3:THR:HG22	44:DN:5:VAL:HG23	1.74	0.70
49:DS:25:ARG:NH2	49:DS:89:ARG:HH12	1.88	0.70
55:DY:74:PRO:O	55:DY:75:ILE:HB	1.89	0.70
1:AA:1191:A:H5''	3:AC:4:LYS:NZ	2.06	0.70
5:AE:12:LEU:HD13	5:AE:31:LEU:HB3	1.72	0.70
5:AE:90:VAL:C	5:AE:91:LEU:HD12	2.11	0.70
7:AG:65:ALA:HB1	7:AG:127:ALA:HB3	1.74	0.70
35:BA:1570:A:H2'	35:BA:1571:A:H8	1.54	0.70
35:BA:2467:C:C2'	35:BA:2468:G:H5'	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:445:C:H5''	51:BU:3:ARG:CB	2.21	0.70
35:BA:713:G:O2'	35:BA:714:U:H5'	1.90	0.70
41:BG:173:LEU:CD2	41:BG:173:LEU:H	2.04	0.70
42:BH:153:LYS:HB2	42:BH:154:PRO:HD2	1.73	0.70
50:BT:22:PHE:CD2	50:BT:22:PHE:N	2.54	0.70
44:BN:42:TRP:N	51:BU:64:ARG:NH1	2.39	0.70
55:BY:28:LYS:HD2	55:BY:37:VAL:CG1	2.21	0.70
55:BY:31:LEU:CD1	55:BY:34:LYS:H	2.03	0.70
56:BZ:107:THR:CG2	56:BZ:111:VAL:HG11	2.22	0.70
1:CA:1346:A:N1	1:CA:1374:A:H5''	2.06	0.70
1:CA:489:C:H2'	1:CA:490:G:C8	2.26	0.70
1:CA:498:U:H2'	1:CA:498:U:O2	1.89	0.70
1:CA:625:G:H2'	1:CA:626:U:C6	2.26	0.70
4:CD:153:ARG:HB3	4:CD:153:ARG:NH1	2.07	0.70
29:D3:29:ARG:HH11	35:DA:1183:G:H4'	1.56	0.70
35:DA:1786:A:N7	35:DA:1938:A:N7	2.40	0.70
35:DA:2007:C:H2'	35:DA:2008:C:H6	1.56	0.70
35:DA:2394:C:OP1	46:DP:63:PRO:HD2	1.89	0.70
35:DA:2672:G:C2'	35:DA:2673:G:H5''	2.21	0.70
45:DO:23:ARG:O	45:DO:39:ILE:HG13	1.91	0.70
50:DT:27:THR:HA	50:DT:88:ILE:H	1.56	0.70
52:DV:3:ALA:HB3	52:DV:14:VAL:HB	1.72	0.70
54:DX:49:VAL:HG12	54:DX:50:LYS:N	2.07	0.70
1:AA:448:A:O2'	1:AA:449:C:H5'	1.91	0.70
1:AA:692:U:H2'	1:AA:694:A:OP2	1.91	0.70
1:AA:956:U:O2'	1:AA:957:U:H5'	1.91	0.70
1:AA:977:A:H2'	1:AA:978:A:H5'	1.73	0.70
10:AJ:30:SER:O	10:AJ:81:THR:HG23	1.90	0.70
25:AY:130:ARG:NH1	25:AY:130:ARG:HG3	2.06	0.70
35:BA:2200:C:H42	35:BA:2223:G:H1	1.39	0.70
35:BA:2579:C:O2'	35:BA:2580:U:H5'	1.90	0.70
35:BA:2732:G:C3'	35:BA:2733:A:H5'	2.22	0.70
35:BA:541:C:H2'	35:BA:542:C:C6	2.26	0.70
35:BA:783:A:H2'	35:BA:784:A:H4'	1.73	0.70
35:BA:1799:G:OP1	38:BD:260:ARG:HD2	1.92	0.70
35:BA:2820:A:H62	39:BE:192:ASN:HB2	1.57	0.70
39:BE:96:PHE:HA	39:BE:100:GLU:OE1	1.91	0.70
46:BP:23:PRO:HD2	46:BP:33:ARG:CZ	2.21	0.70
47:BQ:30:GLY:HA3	47:BQ:107:ALA:HB2	1.72	0.70
48:BR:41:ALA:HB1	48:BR:114:VAL:CG2	2.21	0.70
55:BY:81:LYS:HD3	55:BY:97:ARG:O	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:114:U:H2'	1:CA:115:G:C8	2.26	0.70
1:CA:203:U:H4'	1:CA:216:G:C2	2.26	0.70
1:CA:254:G:O2'	1:CA:255:G:H5'	1.90	0.70
4:CD:119:GLN:HG3	4:CD:123:HIS:HD2	1.55	0.70
7:CG:137:LYS:O	7:CG:141:VAL:HG23	1.91	0.70
10:CJ:82:ILE:O	10:CJ:86:MET:HB3	1.90	0.70
11:CK:59:TYR:CZ	11:CK:63:LEU:HD11	2.25	0.70
35:DA:2114:A:H2'	35:DA:2115:G:H5'	1.73	0.70
28:D2:48:HIS:NE2	35:DA:75:G:H4'	2.06	0.70
35:DA:78:A:H2'	35:DA:79:G:H8	1.54	0.70
38:DD:94:LEU:HD13	38:DD:94:LEU:O	1.91	0.70
39:DE:59:VAL:CG2	39:DE:63:LEU:HA	2.22	0.70
47:DQ:65:PHE:O	47:DQ:66:ILE:HG23	1.91	0.70
49:DS:92:TYR:CD1	49:DS:93:LYS:N	2.59	0.70
1:AA:1249:C:H5''	9:AI:70:LYS:HE2	1.74	0.70
1:AA:429:U:H1'	1:AA:430:A:H5''	1.72	0.70
1:AA:946:A:H2'	1:AA:947:G:C8	2.26	0.70
8:AH:32:LYS:O	8:AH:35:ILE:HG12	1.90	0.70
1:AA:303:A:OP1	12:AL:17:LYS:HE3	1.90	0.70
27:B1:46:LEU:N	27:B1:46:LEU:HD12	2.06	0.70
27:B1:11:ARG:HH11	27:B1:60:PHE:HA	1.56	0.70
34:B8:23:VAL:HG12	34:B8:46:ARG:NH1	2.06	0.70
35:BA:251:A:H5''	46:BP:51:PHE:CE1	2.26	0.70
38:BD:131:LEU:CD1	38:BD:136:ILE:HG12	2.20	0.70
39:BE:116:VAL:HG22	39:BE:117:MET:N	2.04	0.70
39:BE:176:ILE:HB	39:BE:181:LEU:HD23	1.72	0.70
40:BF:198:ALA:O	40:BF:201:VAL:HG12	1.92	0.70
41:BG:137:GLU:HB3	41:BG:140:ILE:HG23	1.74	0.70
42:BH:144:VAL:HG12	42:BH:148:ILE:HD11	1.74	0.70
42:BH:70:THR:O	42:BH:72:ILE:N	2.24	0.70
46:BP:79:ARG:HH21	46:BP:109:GLY:CA	2.05	0.70
48:BR:61:HIS:O	48:BR:65:LEU:HB2	1.92	0.70
50:BT:109:GLU:HB3	50:BT:113:LYS:CE	2.18	0.70
53:BW:13:SER:HB3	53:BW:16:LYS:HD2	1.72	0.70
56:BZ:17:ALA:HA	56:BZ:20:ARG:CB	2.21	0.70
1:CA:764:C:H2'	1:CA:765:G:C8	2.26	0.70
5:CE:35:GLY:HA3	5:CE:41:VAL:HG12	1.74	0.70
12:CL:75:HIS:HD2	12:CL:77:LEU:HG	1.55	0.70
13:CM:91:ARG:HB3	13:CM:98:VAL:HG22	1.74	0.70
27:D1:41:ARG:NH1	27:D1:41:ARG:HG3	2.05	0.70
27:D1:89:GLU:OE2	27:D1:90:ILE:HG12	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D2:29:LYS:HG2	28:D2:32:LEU:HD23	1.72	0.70
35:DA:1947:C:H2'	35:DA:1948:G:C8	2.27	0.70
35:DA:49:A:OP1	35:DA:51:G:H5'	1.92	0.70
35:DA:986:C:O2'	35:DA:987:G:H5'	1.91	0.70
38:DD:117:VAL:HG22	38:DD:118:VAL:N	2.06	0.70
38:DD:158:ALA:O	38:DD:161:THR:HG23	1.92	0.70
40:DF:160:ASN:ND2	40:DF:162:LEU:H	1.89	0.70
40:DF:18:ARG:HG2	40:DF:19:GLU:H	1.56	0.70
41:DG:35:GLU:HA	41:DG:99:MET:HE1	1.74	0.70
43:DI:133:HIS:O	43:DI:135:GLU:HG3	1.92	0.70
48:DR:116:LEU:O	48:DR:117:VAL:HB	1.92	0.70
49:DS:34:HIS:CD2	49:DS:53:SER:HB3	2.27	0.70
50:DT:28:VAL:HG11	50:DT:46:GLU:HA	1.74	0.70
44:DN:41:ASP:C	51:DU:64:ARG:HH11	1.95	0.70
53:DW:51:LEU:HD13	53:DW:52:GLU:N	2.07	0.70
54:DX:18:TYR:HA	54:DX:21:PHE:CD1	2.25	0.70
55:DY:81:LYS:HD3	55:DY:97:ARG:O	1.91	0.70
1:AA:1350:A:OP2	9:AI:118:LYS:HD2	1.92	0.70
1:AA:1369:C:OP2	9:AI:111:ARG:HA	1.92	0.70
1:AA:203:U:H4'	1:AA:216:G:C2	2.27	0.70
1:AA:490:G:H2'	1:AA:491:G:H8	1.56	0.70
11:AK:103:LEU:CD2	11:AK:103:LEU:H	2.00	0.70
6:AF:62:TRP:HB2	18:AR:35:ARG:HH12	1.55	0.70
35:BA:2036:C:H5'	35:BA:2036:C:C6	2.19	0.70
39:BE:11:MET:CB	39:BE:24:THR:HA	2.21	0.70
39:BE:179:GLU:HB3	39:BE:181:LEU:CD2	2.20	0.70
39:BE:36:ARG:HH22	39:BE:88:GLY:CA	2.05	0.70
43:BI:110:ASP:O	43:BI:112:LYS:N	2.24	0.70
46:BP:23:PRO:HD2	46:BP:33:ARG:NH1	2.06	0.70
46:BP:74:GLU:OE2	46:BP:75:ILE:HD12	1.91	0.70
50:BT:38:ASN:HD22	50:BT:40:THR:N	1.90	0.70
54:BX:58:HIS:O	54:BX:59:VAL:HG13	1.91	0.70
1:CA:562:C:H1'	12:CL:15:ARG:HB3	1.74	0.70
25:CY:26:ALA:O	25:CY:37:LEU:HA	1.92	0.70
35:DA:1249:U:H5'	35:DA:1249:U:H6	1.57	0.70
35:DA:1791:A:O3'	38:DD:206:LEU:HB2	1.91	0.70
41:DG:132:ASN:HB2	41:DG:159:VAL:CG2	2.20	0.70
45:DO:63:VAL:CG2	45:DO:84:ALA:HA	2.21	0.70
47:DQ:30:GLY:HA3	47:DQ:107:ALA:HB2	1.72	0.70
56:DZ:152:ALA:N	56:DZ:169:GLU:O	2.25	0.70
1:AA:1375:A:H2'	1:AA:1376:U:C6	2.27	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:71:PRO:HG3	7:AG:103:TRP:HZ3	1.54	0.70
10:AJ:5:ARG:HH21	10:AJ:99:LYS:HG3	1.56	0.70
17:AQ:40:LYS:HD2	17:AQ:42:TYR:CE1	2.27	0.70
1:AA:693:G:N2	23:AW:38:A:H2	1.88	0.70
35:BA:1031:G:H22	35:BA:1124:C:H1'	1.57	0.70
35:BA:1899:G:N2	35:BA:1902:C:N4	2.39	0.70
35:BA:2327:A:H2'	35:BA:2328:A:C8	2.26	0.70
35:BA:2852:G:H2'	35:BA:2853:C:C6	2.26	0.70
35:BA:571:A:C5'	35:BA:2030:A:H62	2.05	0.70
35:BA:582:G:H2'	35:BA:583:G:C8	2.25	0.70
45:BO:43:VAL:HG21	45:BO:52:VAL:HG12	1.74	0.70
46:BP:131:SER:HB2	46:BP:134:ALA:CB	2.21	0.70
55:BY:37:VAL:HG23	55:BY:38:ILE:N	2.04	0.70
55:BY:50:ARG:HB2	55:BY:53:PRO:HG3	1.74	0.70
1:CA:1410:G:H2'	1:CA:1411:C:H6	1.56	0.70
1:CA:1432:G:OP1	50:DT:107:ASP:HB2	1.91	0.70
1:CA:1495:U:H2'	1:CA:1496:C:C6	2.27	0.70
1:CA:6:G:H4'	1:CA:298:A:H4'	1.73	0.70
1:CA:336:C:O2'	1:CA:337:C:H5'	1.90	0.70
8:CH:6:ILE:HG22	8:CH:10:LEU:HD11	1.72	0.70
11:CK:69:ALA:HA	11:CK:72:ALA:HB3	1.74	0.70
13:CM:52:GLU:O	13:CM:56:LEU:HB2	1.92	0.70
36:DB:104:U:O2'	36:DB:105:A:H5'	1.91	0.70
38:DD:4:LYS:NZ	38:DD:20:ASP:HA	2.07	0.70
39:DE:49:LEU:N	39:DE:49:LEU:HD22	2.07	0.70
40:DF:8:GLN:HB2	40:DF:126:VAL:HA	1.74	0.70
41:DG:60:LEU:HA	41:DG:63:ILE:HG13	1.74	0.70
42:DH:103:LEU:HD23	42:DH:115:VAL:HB	1.72	0.70
47:DQ:52:VAL:CG1	47:DQ:53:ALA:H	1.99	0.70
51:DU:31:SER:HB3	51:DU:34:LYS:HB2	1.74	0.70
52:DV:39:LEU:HD11	52:DV:53:GLU:N	2.06	0.70
55:DY:28:LYS:HD2	55:DY:37:VAL:HG12	1.72	0.70
1:AA:1507:A:C2	1:AA:1530:G:H1'	2.27	0.70
1:AA:332:G:H2'	1:AA:333:G:H8	1.56	0.70
1:AA:376:G:H5''	16:AP:5:ARG:HD2	1.73	0.70
4:AD:117:ALA:O	4:AD:121:VAL:HG23	1.92	0.70
12:AL:102:ARG:HH11	12:AL:102:ARG:HG2	1.56	0.70
12:AL:86:ARG:HG2	12:AL:87:GLY:H	1.55	0.70
12:AL:85:ILE:HD11	12:AL:98:TYR:HB3	1.72	0.70
1:AA:658:G:C1'	15:AO:22:THR:HB	2.21	0.70
17:AQ:45:HIS:HB2	17:AQ:69:LYS:HE2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:AY:59:THR:O	25:AY:67:VAL:HG22	1.91	0.70
28:B2:57:ILE:HG12	28:B2:59:ARG:NH1	2.06	0.70
31:B5:40:LYS:HE2	31:B5:46:CYS:HB3	1.74	0.70
34:B8:52:LYS:HE3	34:B8:52:LYS:HA	1.73	0.70
35:BA:364:C:C2'	35:BA:365:C:H5''	2.22	0.70
36:BB:11:C:H3'	36:BB:12:C:H6	1.56	0.70
39:BE:173:VAL:HG12	39:BE:174:ASP:H	1.57	0.70
40:BF:160:ASN:ND2	40:BF:162:LEU:H	1.89	0.70
41:BG:64:THR:HG23	41:BG:65:GLY:H	1.57	0.70
43:BI:62:LYS:O	43:BI:62:LYS:HD3	1.91	0.70
43:BI:88:ILE:HG22	43:BI:89:TYR:N	2.06	0.70
47:BQ:64:ILE:HG23	47:BQ:106:VAL:HG13	1.74	0.70
47:BQ:76:LYS:H	47:BQ:88:GLY:CA	2.04	0.70
54:BX:29:TRP:HZ3	54:BX:76:ARG:HG2	1.57	0.70
56:BZ:58:VAL:HG22	56:BZ:68:PRO:CA	2.22	0.70
1:CA:1203:C:H2'	1:CA:1204:A:H8	1.57	0.70
1:CA:471:G:H2'	1:CA:472:A:H8	1.56	0.70
1:CA:924:C:H2'	1:CA:925:G:H8	1.57	0.70
1:CA:974:A:H1'	14:CN:31:ARG:HH21	1.56	0.70
2:CB:101:MET:HA	2:CB:108:ILE:HG13	1.74	0.70
4:CD:194:LEU:N	4:CD:194:LEU:HD22	2.07	0.70
5:CE:149:GLU:O	5:CE:153:LYS:HG2	1.92	0.70
6:CF:33:TYR:CD1	6:CF:75:LEU:HG	2.26	0.70
1:CA:972:C:O3'	10:CJ:57:LYS:HG2	1.91	0.70
17:CQ:82:MET:O	17:CQ:85:VAL:HB	1.92	0.70
25:CY:64:ARG:HA	25:CY:103:ILE:CD1	2.21	0.70
28:D2:48:HIS:CD2	35:DA:75:G:HO2'	2.10	0.70
33:D7:48:LYS:N	33:D7:48:LYS:HD3	2.06	0.70
35:DA:189:G:H2'	35:DA:205:G:N2	2.07	0.70
35:DA:2279:G:N2	35:DA:2280:G:H1'	2.06	0.70
35:DA:2720:U:H5'	35:DA:2721:A:OP2	1.90	0.70
35:DA:389:G:N1	46:DP:71:VAL:HB	2.04	0.70
35:DA:661:C:H2'	35:DA:662:G:C8	2.27	0.70
39:DE:78:LEU:CD2	39:DE:78:LEU:H	2.05	0.70
45:DO:107:ARG:NH1	50:DT:36:GLU:H	1.89	0.70
46:DP:122:PRO:HB3	46:DP:141:ALA:HB1	1.73	0.70
50:DT:83:ILE:HG13	50:DT:84:GLN:HG2	1.72	0.70
52:DV:70:ILE:HB	52:DV:90:PRO:CB	2.19	0.70
1:AA:243:A:H4'	1:AA:244:U:O5'	1.90	0.70
1:AA:254:G:O2'	1:AA:255:G:H5'	1.92	0.70
1:AA:986:A:H1'	19:AS:54:GLY:O	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:19:HIS:O	2:AB:39:ILE:HG23	1.91	0.70
4:AD:59:ARG:HH22	4:AD:66:ARG:NH2	1.89	0.70
10:AJ:56:HIS:O	10:AJ:58:ASP:N	2.25	0.70
17:AQ:9:VAL:HG12	17:AQ:10:VAL:H	1.57	0.70
35:BA:1301:A:H4'	35:BA:1302:A:OP1	1.91	0.70
35:BA:1777:U:O2'	35:BA:1778:U:H5'	1.91	0.70
35:BA:185:U:H2'	35:BA:186:G:C8	2.27	0.70
35:BA:2051:A:H4'	39:BE:141:ILE:CD1	2.22	0.70
35:BA:2443:C:O2'	35:BA:2444:G:H5'	1.92	0.70
35:BA:2506:U:H4'	35:BA:2507:C:OP1	1.91	0.70
35:BA:2801(A):A:H4'	35:BA:2802:G:H5'	1.74	0.70
1:CA:1366:C:H2'	1:CA:1367:C:C6	2.26	0.70
1:CA:137:C:H42	1:CA:226:G:H1	1.37	0.70
1:CA:1513:A:H2'	1:CA:1514:C:C6	2.27	0.70
4:CD:128:VAL:C	4:CD:130:GLY:H	1.93	0.70
4:CD:13:ARG:HG2	4:CD:14:ARG:N	2.07	0.70
4:CD:18:LYS:NZ	4:CD:31:CYS:SG	2.65	0.70
6:CF:39:LYS:HG2	6:CF:40:VAL:H	1.57	0.70
1:CA:1377:A:H2'	7:CG:7:ALA:HB3	1.72	0.70
25:CY:147:LEU:HD23	25:CY:148:HIS:N	2.06	0.70
35:DA:1493:C:H4'	35:DA:1494:A:OP1	1.91	0.70
35:DA:2186:G:C2'	35:DA:2187:G:H5''	2.22	0.70
35:DA:660:G:H5'	40:DF:99:TYR:CD2	2.26	0.70
35:DA:877:U:C2'	35:DA:878:A:H5''	2.22	0.70
40:DF:32:LEU:C	40:DF:32:LEU:HD23	2.11	0.70
41:DG:72:ARG:HD3	41:DG:86:MET:HA	1.72	0.70
44:DN:14:VAL:HG12	44:DN:15:LEU:N	2.03	0.70
52:DV:61:VAL:HG23	52:DV:100:ARG:H	1.55	0.70
55:DY:2:ARG:N	55:DY:4:LYS:HE2	2.06	0.70
1:AA:1288:A:H1'	1:AA:1352:C:O2'	1.92	0.69
1:AA:1507:A:H2'	1:AA:1508:G:H8	1.56	0.69
1:AA:1206:G:H4'	3:AC:192:THR:O	1.92	0.69
1:AA:750:G:N3	15:AO:23:GLY:HA3	2.07	0.69
1:AA:235:C:H1'	17:AQ:61:GLU:OE1	1.92	0.69
17:AQ:95:TYR:O	17:AQ:97:SER:N	2.25	0.69
27:B1:19:GLN:NE2	35:BA:379:G:H21	1.90	0.69
35:BA:1844:C:O2'	35:BA:1845:G:H5'	1.92	0.69
35:BA:2125:G:N2	35:BA:2173:A:H62	1.85	0.69
35:BA:2282:G:H1	35:BA:2427:C:N4	1.89	0.69
35:BA:588:U:H2'	35:BA:589:C:H6	1.55	0.69
35:BA:897:C:H1'	35:BA:899:A:N7	2.05	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:79:VAL:HG12	38:BD:113:VAL:HA	1.73	0.69
41:BG:83:ARG:HB3	41:BG:84:LYS:HD2	1.74	0.69
43:BI:115:ALA:N	43:BI:131:LYS:HE2	2.07	0.69
44:BN:22:THR:CA	44:BN:61:ARG:HB2	2.20	0.69
50:BT:30:VAL:HG12	50:BT:44:ASP:CA	2.22	0.69
50:BT:64:ARG:HB2	50:BT:73:GLU:HB3	1.73	0.69
52:BV:70:ILE:HB	52:BV:90:PRO:CB	2.19	0.69
35:BA:58:G:OP1	54:BX:72:LYS:HB3	1.91	0.69
56:BZ:142:SER:H	56:BZ:144:LEU:CD2	2.05	0.69
1:CA:1190:G:OP1	3:CC:5:ILE:HD12	1.92	0.69
1:CA:688:G:H2'	1:CA:689:C:H6	1.55	0.69
3:CC:124:ILE:HG13	3:CC:130:VAL:HG22	1.74	0.69
5:CE:82:VAL:HG21	5:CE:138:ALA:CA	2.22	0.69
8:CH:45:ILE:HA	8:CH:64:LYS:HB3	1.73	0.69
10:CJ:5:ARG:HH21	10:CJ:99:LYS:HG3	1.57	0.69
12:CL:76:ASN:OD1	12:CL:108:ALA:HB3	1.90	0.69
13:CM:66:LEU:CA	13:CM:70:LEU:HD12	2.22	0.69
18:CR:81:PHE:O	18:CR:82:THR:HB	1.92	0.69
19:CS:53:ASN:HD21	19:CS:56:GLN:N	1.89	0.69
23:CW:57:C:H2'	23:CW:58:A:H8	1.57	0.69
35:DA:1345:C:O2'	35:DA:1346:G:H5'	1.92	0.69
35:DA:1789:A:H2'	35:DA:1790:C:H6	1.56	0.69
35:DA:1999:C:H2'	35:DA:2000:G:H8	1.57	0.69
35:DA:2203:U:O4'	38:DD:151:LYS:HE3	1.92	0.69
35:DA:2312:U:H2'	35:DA:2313:C:H5''	1.74	0.69
35:DA:2801(A):A:H4'	35:DA:2802:G:C2'	2.22	0.69
40:DF:198:ALA:O	40:DF:201:VAL:HG12	1.90	0.69
44:DN:62:VAL:O	44:DN:63:THR:HG22	1.91	0.69
1:CA:1422:G:H4'	45:DO:49:ARG:NH1	2.07	0.69
47:DQ:116:GLU:O	47:DQ:119:ARG:HB3	1.92	0.69
48:DR:17:ARG:HH11	48:DR:17:ARG:HG2	1.57	0.69
50:DT:25:GLY:HA2	50:DT:92:GLY:N	2.06	0.69
52:DV:32:THR:HG22	52:DV:33:VAL:N	2.05	0.69
54:DX:57:LEU:HD12	54:DX:76:ARG:NE	2.06	0.69
1:AA:349:A:O2'	1:AA:350:G:H5'	1.91	0.69
2:AB:233:SER:HB2	2:AB:234:PRO:HD2	1.72	0.69
17:AQ:11:VAL:HA	17:AQ:53:LEU:HD11	1.73	0.69
36:BB:79:C:O2'	36:BB:80:U:H5'	1.91	0.69
40:BF:37:VAL:O	40:BF:40:GLN:HB2	1.93	0.69
41:BG:101:ILE:CD1	41:BG:105:LYS:HE3	2.22	0.69
41:BG:19:LEU:HD21	41:BG:175:LEU:CD1	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BQ:10:ARG:HD3	47:BQ:12:GLN:HB3	1.75	0.69
53:BW:70:TYR:HE2	53:BW:108:GLY:HA3	1.56	0.69
54:BX:81:VAL:HG13	54:BX:85:PRO:HB2	1.73	0.69
1:CA:1456:G:H2'	1:CA:1457:G:C5'	2.22	0.69
2:CB:102:LEU:CD1	2:CB:102:LEU:H	2.04	0.69
2:CB:54:THR:HG22	2:CB:58:ILE:HD11	1.74	0.69
2:CB:69:LEU:CD1	2:CB:71:VAL:HG23	2.22	0.69
6:CF:48:LEU:HD22	18:CR:77:GLY:HA3	1.72	0.69
6:CF:53:ALA:HB3	6:CF:86:ARG:NH1	2.05	0.69
1:CA:668:G:O4'	15:CO:49:ASP:HB2	1.92	0.69
19:CS:62:ILE:HD12	19:CS:63:THR:H	1.56	0.69
20:CT:57:ARG:NH1	20:CT:57:ARG:HB2	2.07	0.69
24:CX:20:A:H2'	24:CX:21:G:C8	2.27	0.69
25:CY:14:MET:O	25:CY:18:LEU:HB3	1.91	0.69
35:DA:1751:C:O2'	35:DA:1752:C:H5'	1.92	0.69
35:DA:1930:G:H22	35:DA:1968:G:H2'	1.58	0.69
41:DG:129:GLY:C	41:DG:131:TYR:H	1.96	0.69
43:DI:113:ARG:HA	43:DI:131:LYS:HE3	1.73	0.69
45:DO:2:ILE:HD11	45:DO:82:ASN:CB	2.22	0.69
46:DP:80:TYR:CE1	46:DP:111:ARG:HB3	2.27	0.69
46:DP:17:LYS:C	46:DP:19:VAL:H	1.95	0.69
47:DQ:66:ILE:HG22	47:DQ:104:PHE:CE2	2.26	0.69
55:DY:28:LYS:HA	55:DY:39:VAL:H	1.56	0.69
5:AE:82:VAL:HG21	5:AE:138:ALA:CA	2.21	0.69
7:AG:120:ILE:HG22	7:AG:124:LEU:HD11	1.73	0.69
11:AK:61:ALA:HB3	11:AK:90:GLY:HA3	1.72	0.69
35:BA:1340:U:C6	35:BA:1603:A:O4'	2.45	0.69
35:BA:1600:C:O2'	35:BA:1601:G:H5'	1.92	0.69
35:BA:709:U:H2'	35:BA:710:G:H8	1.56	0.69
35:BA:7:G:H4'	44:BN:13:TRP:CH2	2.27	0.69
35:BA:869:G:H1'	47:BQ:8:LYS:NZ	2.06	0.69
36:BB:104:U:O2'	36:BB:105:A:H5'	1.92	0.69
41:BG:166:ASP:OD1	41:BG:170:ARG:HB2	1.92	0.69
47:BQ:132:VAL:HG11	56:BZ:81:ARG:NH1	2.07	0.69
47:BQ:132:VAL:HG12	47:BQ:133:ARG:H	1.56	0.69
55:BY:76:CYS:O	55:BY:78:ALA:N	2.25	0.69
56:BZ:76:LEU:O	56:BZ:78:LYS:N	2.24	0.69
1:CA:1203:C:OP1	14:CN:3:ARG:HD2	1.92	0.69
1:CA:522:C:H41	12:CL:53:ARG:NH2	1.86	0.69
10:CJ:5:ARG:HA	10:CJ:73:ASP:OD1	1.93	0.69
1:CA:750:G:N3	15:CO:23:GLY:HA3	2.06	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:47:THR:O	18:CR:82:THR:HA	1.92	0.69
18:CR:67:ALA:HA	18:CR:70:ILE:HD12	1.73	0.69
19:CS:12:ASP:HB3	19:CS:15:LEU:HD23	1.75	0.69
35:DA:1193:G:H2'	35:DA:1194:A:O4'	1.92	0.69
35:DA:1258:C:H2'	35:DA:1259:G:C8	2.28	0.69
35:DA:185:U:H2'	35:DA:186:G:C8	2.27	0.69
35:DA:2467:C:C2'	35:DA:2468:G:H5'	2.23	0.69
35:DA:2845:G:O2'	35:DA:2846:G:H5'	1.92	0.69
35:DA:445:C:H5''	51:DU:3:ARG:CB	2.22	0.69
37:DC:58:VAL:CG2	37:DC:166:ASP:H	2.04	0.69
38:DD:79:VAL:CG1	38:DD:113:VAL:HA	2.22	0.69
39:DE:36:ARG:HH22	39:DE:88:GLY:CA	2.05	0.69
43:DI:68:LEU:O	43:DI:68:LEU:HD23	1.92	0.69
44:DN:120:LEU:HD11	44:DN:122:VAL:CG2	2.23	0.69
47:DQ:8:LYS:CG	47:DQ:9:TYR:H	2.04	0.69
48:DR:4:LEU:O	48:DR:4:LEU:HD13	1.92	0.69
52:DV:19:LYS:HG3	52:DV:20:LEU:N	2.07	0.69
1:AA:489:C:H2'	1:AA:490:G:C8	2.24	0.69
4:AD:96:LEU:N	4:AD:96:LEU:HD22	2.02	0.69
6:AF:21:LEU:O	6:AF:24:GLU:HG2	1.93	0.69
6:AF:53:ALA:HB3	6:AF:86:ARG:NH1	2.08	0.69
16:AP:6:LEU:HB3	16:AP:17:TYR:HD2	1.58	0.69
35:BA:1177:A:H5'	35:BA:1178:C:C5	2.28	0.69
35:BA:137:C:H2'	35:BA:139:G:H5'	1.74	0.69
35:BA:186:G:O2'	35:BA:187:G:H5'	1.92	0.69
35:BA:2312:U:H2'	35:BA:2313:C:H5''	1.73	0.69
39:BE:116:VAL:O	39:BE:117:MET:HB3	1.91	0.69
40:BF:129:PHE:HA	40:BF:142:TRP:NE1	2.06	0.69
52:BV:61:VAL:CG2	52:BV:100:ARG:H	2.05	0.69
53:BW:51:LEU:C	53:BW:51:LEU:HD13	2.13	0.69
54:BX:49:VAL:HG12	54:BX:50:LYS:N	2.07	0.69
56:BZ:165:VAL:HG12	56:BZ:166:SER:N	2.07	0.69
1:CA:1292:U:H2'	1:CA:1293:G:C8	2.28	0.69
1:CA:1320:C:N4	19:CS:36:ARG:HG3	2.08	0.69
1:CA:1324:A:H2'	1:CA:1325:C:C6	2.19	0.69
1:CA:444:C:H2'	1:CA:445:G:H8	1.55	0.69
1:CA:955:U:H1'	1:CA:1227:A:N6	2.07	0.69
6:CF:21:LEU:O	6:CF:24:GLU:HG2	1.93	0.69
10:CJ:45:ARG:O	10:CJ:64:GLU:HA	1.93	0.69
27:D1:16:ASN:HD22	27:D1:16:ASN:H	1.39	0.69
35:DA:1177:A:H5'	35:DA:1178:C:C5	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1341:U:P	35:DA:1397:U:H3	2.15	0.69
35:DA:1786:A:C5	35:DA:1938:A:N7	2.61	0.69
35:DA:814:C:O2'	35:DA:815:C:H5'	1.92	0.69
41:DG:40:ASN:ND2	41:DG:91:ARG:HB2	2.08	0.69
46:DP:39:LYS:HD3	46:DP:40:SER:H	1.56	0.69
46:DP:66:GLY:O	46:DP:68:GLN:HG2	1.92	0.69
1:AA:1062:U:H2'	1:AA:1063:C:C6	2.28	0.69
1:AA:1084:G:H5'	1:AA:1102:A:OP2	1.92	0.69
1:AA:439:A:C2	1:AA:441:A:H1'	2.27	0.69
1:AA:724:G:O2'	1:AA:725:G:H5'	1.92	0.69
5:AE:72:GLN:HE22	5:AE:77:PRO:HD3	1.57	0.69
9:AI:5:TYR:HE1	9:AI:7:THR:HG1	1.38	0.69
35:BA:2036:C:H6	35:BA:2036:C:C5'	2.05	0.69
35:BA:2284:C:H2'	35:BA:2285:C:C5'	2.16	0.69
35:BA:2514:U:H2'	35:BA:2515:C:C6	2.27	0.69
35:BA:2701:C:H3'	35:BA:2702:U:C5'	2.20	0.69
35:BA:2820:A:H62	39:BE:192:ASN:CB	2.06	0.69
44:BN:42:TRP:CD1	51:BU:63:VAL:HG11	2.27	0.69
46:BP:47:ASP:OD1	46:BP:49:ARG:HB3	1.92	0.69
55:BY:101:LYS:HG2	55:BY:102:CYS:N	2.05	0.69
1:CA:320:C:H2'	1:CA:321:A:C8	2.27	0.69
3:CC:149:ALA:HA	3:CC:201:TYR:O	1.92	0.69
5:CE:144:THR:O	5:CE:148:VAL:HG23	1.92	0.69
7:CG:71:PRO:HG3	7:CG:103:TRP:HZ3	1.56	0.69
11:CK:58:PRO:HD3	11:CK:89:ALA:HB1	1.75	0.69
25:CY:73:GLN:HB2	25:CY:77:LYS:HZ1	1.56	0.69
31:D5:13:LYS:O	31:D5:16:ARG:HB3	1.93	0.69
32:D6:32:ASN:CG	32:D6:33:LYS:H	1.94	0.69
35:DA:1221:C:H2'	35:DA:1221(A):C:H6	1.57	0.69
35:DA:1509(B):A:H2'	35:DA:1510:G:C8	2.27	0.69
35:DA:2472:G:H5'	35:DA:2473:U:H5''	1.75	0.69
35:DA:575:A:C2'	35:DA:576:U:H5'	2.23	0.69
35:DA:588:U:H2'	35:DA:589:C:H6	1.53	0.69
35:DA:661:C:H4'	46:DP:18:ARG:HG2	1.74	0.69
35:DA:713:G:O2'	35:DA:714:U:H5'	1.93	0.69
35:DA:755:C:H2'	35:DA:756:C:C6	2.26	0.69
44:DN:9:VAL:HG11	44:DN:39:ARG:NH2	2.04	0.69
44:DN:42:TRP:CD1	51:DU:63:VAL:HG11	2.28	0.69
46:DP:17:LYS:O	46:DP:19:VAL:N	2.25	0.69
47:DQ:51:ARG:O	47:DQ:55:VAL:HG13	1.92	0.69
50:DT:78:LEU:HD23	50:DT:79:HIS:CE1	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:104:PHE:HE2	56:DZ:122:ARG:HA	1.57	0.69
1:AA:1287:A:H2'	1:AA:1288:A:C8	2.27	0.69
1:AA:1291:G:H4'	9:AI:39:GLY:HA3	1.74	0.69
1:AA:1402:C:H2'	1:AA:1403:C:O4'	1.92	0.69
1:AA:449:C:O2	16:AP:42:ARG:HD2	1.92	0.69
1:AA:570:G:H2'	1:AA:571:U:C6	2.27	0.69
1:AA:677:U:H3	1:AA:713:G:H22	1.41	0.69
1:AA:818:G:H3'	1:AA:819:A:H5''	1.75	0.69
2:AB:162:ILE:O	2:AB:162:ILE:HD12	1.93	0.69
3:AC:73:PRO:HA	3:AC:76:VAL:HG13	1.73	0.69
10:AJ:45:ARG:O	10:AJ:64:GLU:HA	1.93	0.69
10:AJ:9:ARG:O	10:AJ:94:VAL:HG13	1.92	0.69
27:B1:87:PRO:HB2	27:B1:91:LYS:NZ	2.05	0.69
35:BA:1405:U:H2'	35:BA:1406:U:C6	2.27	0.69
35:BA:2127:G:H1'	35:BA:2128:C:H4'	1.74	0.69
35:BA:2712:U:H1'	35:BA:2712(A):A:C8	2.26	0.69
35:BA:271(U):G:H2'	35:BA:271(V):G:C8	2.27	0.69
35:BA:755:C:H2'	35:BA:756:C:C6	2.27	0.69
38:BD:210:GLY:O	38:BD:211:ARG:HB3	1.90	0.69
42:BH:43:VAL:O	42:BH:43:VAL:HG23	1.92	0.69
43:BI:12:LEU:HD12	43:BI:19:VAL:HG11	1.74	0.69
43:BI:38:LEU:HB2	43:BI:40:THR:HG23	1.75	0.69
47:BQ:50:ALA:HA	47:BQ:124:LYS:HG3	1.75	0.69
47:BQ:140:ALA:HB3	56:BZ:53:ILE:CD1	2.16	0.69
48:BR:63:ARG:O	48:BR:67:LEU:HD23	1.92	0.69
35:BA:2013:A:H4'	53:BW:96:ILE:HD12	1.73	0.69
1:CA:1375:A:H2'	1:CA:1376:U:H6	1.57	0.69
1:CA:689:C:P	11:CK:46:GLY:HA3	2.33	0.69
1:CA:977:A:C2'	1:CA:978:A:H5'	2.22	0.69
3:CC:16:ARG:NH1	3:CC:16:ARG:HB2	2.08	0.69
5:CE:147:ASP:HA	5:CE:150:ARG:HB3	1.73	0.69
25:CY:45:TYR:O	25:CY:46:TYR:HB2	1.92	0.69
34:D8:52:LYS:H	34:D8:53:PRO:CD	2.03	0.69
35:DA:251:A:H5''	46:DP:51:PHE:CE1	2.28	0.69
35:DA:797:C:H2'	35:DA:798:G:C8	2.27	0.69
40:DF:53:THR:CG2	40:DF:56:GLU:HB2	2.22	0.69
41:DG:51:ARG:NE	41:DG:51:ARG:CA	2.45	0.69
41:DG:63:ILE:HD12	41:DG:64:THR:N	2.07	0.69
43:DI:83:ALA:HB1	43:DI:88:ILE:HG12	1.73	0.69
1:AA:696:A:H2'	1:AA:697:U:H6	1.56	0.69
4:AD:142:PRO:HA	4:AD:185:PHE:CD2	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:8:ASN:ND2	17:AQ:34:LYS:HE2	2.08	0.69
20:AT:43:LEU:HD12	20:AT:52:ALA:HA	1.75	0.69
32:B6:15:GLU:O	32:B6:15:GLU:HG2	1.93	0.69
35:BA:2562:U:H2'	35:BA:2563:U:H5'	1.75	0.69
35:BA:575:A:C2'	35:BA:576:U:H5'	2.22	0.69
35:BA:852:G:O2'	35:BA:853:G:H5'	1.93	0.69
37:BC:58:VAL:CG2	37:BC:166:ASP:H	2.05	0.69
38:BD:118:VAL:HG22	38:BD:119:ALA:N	2.07	0.69
38:BD:161:THR:O	38:BD:196:VAL:HG23	1.92	0.69
38:BD:264:LYS:HE2	38:BD:266:SER:O	1.92	0.69
40:BF:9:ILE:HG12	40:BF:14:PRO:HA	1.74	0.69
35:BA:660:G:H5'	40:BF:99:TYR:CE2	2.27	0.69
46:BP:30:THR:CG2	46:BP:31:ALA:H	2.05	0.69
1:CA:1126:U:H2'	1:CA:1127:G:C8	2.27	0.69
1:CA:783:C:O2'	1:CA:784:C:H5'	1.91	0.69
25:CY:30:THR:HG22	25:CY:179:LYS:HE3	1.74	0.69
25:CY:18:LEU:HG	25:CY:19:GLU:HG3	1.75	0.69
35:DA:1232:G:H2'	35:DA:1233:C:H6	1.57	0.69
35:DA:1991:U:H2'	35:DA:1992:G:C5'	2.23	0.69
35:DA:2200:C:H42	35:DA:2223:G:H1	1.38	0.69
35:DA:2723:C:C2'	35:DA:2724:C:H5'	2.23	0.69
35:DA:272(D):G:H1	35:DA:364:C:N4	1.88	0.69
35:DA:2808:U:H2'	35:DA:2809:A:H5'	1.72	0.69
35:DA:296:C:O2'	35:DA:297:C:H5'	1.93	0.69
35:DA:686:G:N2	35:DA:788:A:H61	1.89	0.69
41:DG:137:GLU:OE2	41:DG:140:ILE:HG23	1.92	0.69
44:DN:107:LEU:HB2	44:DN:108:PRO:HD2	1.75	0.69
46:DP:16:ARG:CD	46:DP:18:ARG:H	2.04	0.69
46:DP:58:THR:C	46:DP:60:MET:H	1.96	0.69
46:DP:59:LEU:HA	46:DP:61:ARG:NE	2.08	0.69
56:DZ:33:LEU:HD11	56:DZ:35:ARG:CG	2.23	0.69
1:AA:17:U:H2'	1:AA:18:C:C6	2.27	0.69
1:AA:714:G:H21	1:AA:777:A:H1'	1.57	0.69
5:AE:42:GLY:HA3	5:AE:66:MET:HG2	1.74	0.69
9:AI:65:VAL:HG22	9:AI:66:ARG:N	2.08	0.69
9:AI:70:LYS:O	9:AI:74:ILE:HG13	1.92	0.69
1:AA:881:G:P	12:AL:12:ARG:NH2	2.66	0.69
26:B0:23:VAL:HG13	26:B0:37:LEU:O	1.92	0.69
28:B2:57:ILE:O	28:B2:57:ILE:HG12	1.90	0.69
29:B3:29:ARG:HH11	35:BA:1183:G:H4'	1.56	0.69
35:BA:2114:A:H2'	35:BA:2115:G:H5'	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2680:C:H2'	35:BA:2681:C:O2	1.92	0.69
38:BD:117:VAL:HG22	38:BD:118:VAL:N	2.07	0.69
39:BE:105:THR:OG1	39:BE:166:THR:HG22	1.92	0.69
47:BQ:28:ALA:CB	47:BQ:67:ARG:HD2	2.22	0.69
48:BR:82:GLU:O	48:BR:85:PRO:HD2	1.92	0.69
50:BT:27:THR:O	50:BT:28:VAL:HB	1.91	0.69
1:CA:533:A:O2'	1:CA:534:U:H5''	1.93	0.69
2:CB:69:LEU:HD23	2:CB:159:PRO:HG2	1.75	0.69
9:CI:83:ARG:O	9:CI:86:VAL:HG12	1.93	0.69
12:CL:36:VAL:O	12:CL:58:VAL:HG13	1.93	0.69
12:CL:86:ARG:HG2	12:CL:87:GLY:N	2.08	0.69
35:DA:1188:U:O2'	35:DA:1189:A:H5'	1.93	0.69
35:DA:1799:G:OP1	38:DD:260:ARG:HD2	1.92	0.69
35:DA:571:A:C5'	35:DA:2030:A:H62	2.05	0.69
35:DA:208:C:H2'	35:DA:209:C:C6	2.28	0.69
35:DA:2801(A):A:H4'	35:DA:2802:G:H5'	1.73	0.69
35:DA:2830:G:H5'	39:DE:58:ARG:HH12	1.57	0.69
27:D1:19:GLN:NE2	35:DA:379:G:H21	1.90	0.69
35:DA:897:C:H1'	35:DA:899:A:N7	2.07	0.69
43:DI:129:THR:HA	43:DI:137:PRO:HA	1.75	0.69
43:DI:62:LYS:O	43:DI:62:LYS:HD3	1.93	0.69
52:DV:61:VAL:CG2	52:DV:100:ARG:H	2.05	0.69
54:DX:16:LYS:HE3	54:DX:16:LYS:HA	1.74	0.69
1:AA:498:U:H2'	1:AA:499:A:H5'	1.75	0.69
6:AF:45:LEU:HD12	6:AF:46:ARG:N	2.08	0.69
9:AI:18:PHE:HB2	9:AI:62:TYR:O	1.93	0.69
1:AA:377:G:OP1	16:AP:3:LYS:HD3	1.92	0.69
35:BA:110:G:O2'	35:BA:111:A:H5'	1.92	0.69
35:BA:1449:A:C2	35:BA:1529:G:H1'	2.26	0.69
35:BA:184:C:H2'	35:BA:185:U:C6	2.28	0.69
35:BA:571:A:H5'	35:BA:2030:A:N6	2.07	0.69
35:BA:2317:C:O2'	35:BA:2318:G:H5'	1.93	0.69
39:BE:101:ARG:HD3	39:BE:169:ASN:ND2	2.07	0.69
43:BI:115:ALA:H	43:BI:131:LYS:HE2	1.57	0.69
50:BT:78:LEU:HD23	50:BT:79:HIS:CE1	2.28	0.69
56:BZ:60:GLU:HA	56:BZ:66:SER:HB3	1.75	0.69
1:CA:179:A:H2'	1:CA:180:U:C6	2.27	0.69
1:CA:920:U:H2'	1:CA:921:U:C6	2.28	0.69
3:CC:148:GLY:CA	3:CC:203:PHE:HB3	2.23	0.69
5:CE:13:ILE:HA	5:CE:29:GLY:O	1.93	0.69
9:CI:28:VAL:HG12	9:CI:29:ASN:N	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:972:C:H4'	10:CJ:57:LYS:HG3	1.73	0.69
11:CK:21:ILE:HD13	11:CK:84:VAL:HG12	1.74	0.69
33:D7:15:THR:HG22	33:D7:16:HIS:CD2	2.28	0.69
35:DA:1270:C:H5''	35:DA:1271:G:H5'	1.74	0.69
35:DA:2250:G:C5	47:DQ:82:ARG:HD2	2.28	0.69
35:DA:64:A:O2'	35:DA:65:C:H5'	1.92	0.69
38:DD:267:SER:O	38:DD:269:PHE:N	2.25	0.69
43:DI:4:ILE:HD13	43:DI:47:LEU:HD22	1.75	0.69
46:DP:47:ASP:OD1	46:DP:49:ARG:HB3	1.93	0.69
47:DQ:39:PRO:HB3	47:DQ:99:PRO:CD	2.12	0.69
50:DT:64:ARG:HB2	50:DT:73:GLU:HB3	1.73	0.69
50:DT:75:ILE:HD12	50:DT:75:ILE:N	2.07	0.69
55:DY:90:LEU:HG	55:DY:91:GLU:HG2	1.75	0.69
6:AF:9:VAL:C	6:AF:10:LEU:HD12	2.13	0.69
10:AJ:27:ALA:CB	10:AJ:85:LEU:HD11	2.23	0.69
27:B1:41:ARG:HH22	35:BA:205:G:H1	1.37	0.69
35:BA:1860:G:H1	35:BA:1882:C:H42	1.41	0.69
35:BA:2223:G:C2'	35:BA:2224:G:H5'	2.22	0.69
35:BA:404:C:C4'	35:BA:405:U:H5'	2.20	0.69
38:BD:267:SER:C	38:BD:269:PHE:N	2.41	0.69
41:BG:96:ARG:O	41:BG:99:MET:HB3	1.93	0.69
42:BH:97:ARG:O	42:BH:125:VAL:HG11	1.93	0.69
42:BH:41:MET:HG3	42:BH:53:GLU:O	1.93	0.69
47:BQ:71:ASP:O	47:BQ:73:PRO:HD3	1.93	0.69
54:BX:25:LYS:NZ	54:BX:87:GLN:N	2.41	0.69
54:BX:54:VAL:HG13	54:BX:78:LYS:O	1.92	0.69
1:CA:1512:U:H2'	1:CA:1513:A:C8	2.28	0.69
1:CA:67:C:H2'	1:CA:68:G:C8	2.28	0.69
8:CH:121:ASP:O	8:CH:125:ARG:HB2	1.93	0.69
35:DA:1405:U:H2'	35:DA:1406:U:C6	2.28	0.69
35:DA:184:C:H2'	35:DA:185:U:C6	2.27	0.69
35:DA:2712:U:H1'	35:DA:2712(A):A:C8	2.28	0.69
38:DD:45:ASN:CG	38:DD:46:GLN:N	2.46	0.69
35:DA:7:G:H4'	44:DN:13:TRP:CH2	2.28	0.69
51:DU:83:LEU:CB	51:DU:88:ILE:HG12	2.23	0.69
55:DY:50:ARG:HB2	55:DY:53:PRO:HG3	1.73	0.69
1:AA:922:G:C2	1:AA:1396:A:C2	2.81	0.69
1:AA:522:C:N4	12:AL:53:ARG:HH21	1.88	0.69
19:AS:12:ASP:HB3	19:AS:15:LEU:HD23	1.75	0.69
19:AS:33:THR:HG23	19:AS:51:VAL:HA	1.75	0.69
25:AY:118:VAL:C	25:AY:120:GLN:H	1.94	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:AY:80:GLU:C	25:AY:82:ALA:H	1.97	0.69
35:BA:1324:G:H3'	35:BA:1325:G:H4'	1.75	0.69
38:BD:101:GLU:OE1	38:BD:103:ARG:HD3	1.93	0.69
38:BD:80:ALA:HB3	38:BD:94:LEU:CD1	2.23	0.69
39:BE:1:MET:HG2	39:BE:83:ASP:O	1.92	0.69
43:BI:77:LEU:CB	43:BI:140:LEU:HD13	2.18	0.69
43:BI:83:ALA:HB1	43:BI:88:ILE:HG12	1.73	0.69
46:BP:101:VAL:CG2	46:BP:107:LYS:HA	2.23	0.69
49:BS:26:LEU:HA	49:BS:39:ILE:HG13	1.74	0.69
50:BT:75:ILE:HD12	50:BT:75:ILE:N	2.08	0.69
1:CA:920:U:H1'	1:CA:1080:A:H2	1.55	0.69
1:CA:976:G:H22	1:CA:1362:C:H2'	1.57	0.69
1:CA:1369:C:OP2	9:CI:111:ARG:HA	1.92	0.69
1:CA:794:A:H4'	1:CA:1521:G:O2'	1.93	0.69
1:CA:126:G:H5'	1:CA:633:G:N2	2.07	0.69
1:CA:714:G:H21	1:CA:777:A:H1'	1.58	0.69
1:CA:818:G:H3'	1:CA:819:A:H5''	1.74	0.69
2:CB:187:LEU:C	2:CB:187:LEU:HD22	2.14	0.69
5:CE:102:ALA:HB2	5:CE:120:THR:OG1	1.92	0.69
6:CF:14:LEU:CD1	6:CF:19:LEU:HB2	2.21	0.69
1:CA:673:G:H5''	6:CF:87:ARG:NH1	2.08	0.69
9:CI:18:PHE:HB2	9:CI:62:TYR:O	1.93	0.69
17:CQ:19:VAL:CG2	17:CQ:44:ALA:HB3	2.23	0.69
35:DA:990:A:N6	35:DA:1186:G:H1'	2.08	0.69
35:DA:1830:C:H42	35:DA:1975:G:H1	1.40	0.69
35:DA:1991:U:H2'	35:DA:1992:G:H5'	1.75	0.69
35:DA:2127:G:H1'	35:DA:2128:C:H4'	1.75	0.69
35:DA:603:A:H4'	35:DA:604:G:O5'	1.92	0.69
35:DA:678:C:H2'	35:DA:679:C:C6	2.25	0.69
35:DA:696:G:C2'	35:DA:697:C:H5'	2.23	0.69
41:DG:114:ILE:HG22	41:DG:115:ARG:N	2.06	0.69
42:DH:145:ALA:HB1	42:DH:164:TYR:HE1	1.58	0.69
53:DW:70:TYR:HE2	53:DW:108:GLY:HA3	1.58	0.69
56:DZ:134:PRO:O	56:DZ:136:PHE:N	2.25	0.69
1:AA:67:C:H2'	1:AA:68:G:C8	2.27	0.68
3:AC:15:THR:CG2	3:AC:16:ARG:HH12	2.00	0.68
5:AE:36:ASP:O	5:AE:37:ARG:HB2	1.92	0.68
5:AE:51:VAL:O	5:AE:55:VAL:HG23	1.93	0.68
8:AH:12:ARG:NH1	8:AH:26:VAL:HA	2.09	0.68
10:AJ:5:ARG:O	10:AJ:98:ILE:HA	1.92	0.68
13:AM:75:ALA:O	13:AM:79:LYS:HG3	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:19:VAL:CG2	17:AQ:44:ALA:HB3	2.22	0.68
18:AR:47:THR:O	18:AR:82:THR:HA	1.92	0.68
18:AR:47:THR:OG1	18:AR:49:LYS:HG2	1.93	0.68
23:AW:19:G:H3'	23:AW:20:G:C5'	2.23	0.68
35:BA:1221(A):C:O2'	35:BA:1222:C:H5'	1.92	0.68
35:BA:1528(A):A:C3'	35:BA:1529:G:H5''	2.22	0.68
35:BA:1300:U:O2	35:BA:1626:G:H2'	1.92	0.68
35:BA:2807:G:H3'	35:BA:2808:U:H5''	1.74	0.68
35:BA:660:G:H5'	40:BF:99:TYR:CD2	2.28	0.68
35:BA:833:U:H2'	35:BA:834:C:C6	2.27	0.68
35:BA:920:G:H2'	35:BA:921:G:H8	1.56	0.68
38:BD:206:LEU:HA	38:BD:211:ARG:HH12	1.57	0.68
38:BD:25:THR:O	38:BD:26:LYS:HD2	1.93	0.68
43:BI:23:PRO:HB3	43:BI:27:ARG:HH21	1.57	0.68
45:BO:68:GLU:CD	45:BO:78:ARG:HH11	1.97	0.68
47:BQ:34:LEU:HD11	47:BQ:129:THR:CB	2.23	0.68
54:BX:25:LYS:HZ3	54:BX:87:GLN:N	1.92	0.68
1:CA:1416:G:H2'	1:CA:1417:G:O4'	1.93	0.68
1:CA:254:G:OP1	17:CQ:67:LYS:O	2.10	0.68
1:CA:962:C:H2'	1:CA:963:G:H8	1.58	0.68
2:CB:84:GLU:CB	2:CB:219:VAL:HG21	2.20	0.68
6:CF:22:GLU:O	6:CF:26:ILE:HG13	1.93	0.68
7:CG:152:ALA:C	7:CG:154:TYR:H	1.97	0.68
14:CN:34:TYR:C	14:CN:36:PHE:H	1.96	0.68
15:CO:56:LEU:O	15:CO:60:VAL:HG23	1.93	0.68
15:CO:36:ILE:HD12	15:CO:63:ARG:HD3	1.76	0.68
32:D6:15:GLU:O	32:D6:15:GLU:HG2	1.93	0.68
33:D7:48:LYS:HD3	33:D7:48:LYS:H	1.58	0.68
35:DA:1258:C:H2'	35:DA:1259:G:H8	1.57	0.68
35:DA:1528(A):A:C3'	35:DA:1529:G:H5''	2.22	0.68
35:DA:2175:C:C2'	35:DA:2176:A:H5''	2.22	0.68
35:DA:226:G:H5'	35:DA:257:A:H4'	1.75	0.68
35:DA:2808:U:C2'	35:DA:2809:A:H5'	2.23	0.68
38:DD:210:GLY:O	38:DD:212:SER:N	2.26	0.68
39:DE:77:ILE:HG23	39:DE:78:LEU:HD23	1.75	0.68
49:DS:38:GLN:HG2	49:DS:39:ILE:N	2.06	0.68
9:AI:28:VAL:HG12	9:AI:29:ASN:N	2.08	0.68
10:AJ:27:ALA:HB2	10:AJ:85:LEU:HD11	1.74	0.68
27:B1:20:ARG:HH12	27:B1:41:ARG:NH2	1.89	0.68
33:B7:48:LYS:H	33:B7:48:LYS:HD3	1.58	0.68
35:BA:1937:A:C2'	35:BA:1938:A:H5'	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1791:A:O3'	38:BD:206:LEU:HB2	1.93	0.68
38:BD:35:LYS:CE	38:BD:104:TYR:HB2	2.22	0.68
39:BE:59:VAL:HG11	39:BE:63:LEU:HG	1.74	0.68
35:BA:2787:C:O2	39:BE:61:ARG:HD3	1.93	0.68
41:BG:172:LEU:O	41:BG:175:LEU:HB2	1.93	0.68
35:BA:1190:G:H4'	46:BP:35:HIS:HB3	1.76	0.68
52:BV:61:VAL:HG23	52:BV:100:ARG:H	1.56	0.68
54:BX:81:VAL:CG1	54:BX:85:PRO:HB2	2.23	0.68
1:CA:1401:G:H2'	1:CA:1402:C:H5'	1.75	0.68
3:CC:109:PRO:HA	3:CC:115:LEU:CD1	2.23	0.68
5:CE:36:ASP:O	5:CE:37:ARG:HB2	1.93	0.68
7:CG:120:ILE:HG22	7:CG:124:LEU:HD11	1.74	0.68
10:CJ:51:ARG:H	10:CJ:60:ARG:HA	1.58	0.68
13:CM:23:TYR:CD1	13:CM:67:GLU:HA	2.27	0.68
17:CQ:11:VAL:HA	17:CQ:53:LEU:HD11	1.74	0.68
19:CS:63:THR:CG2	19:CS:66:MET:HG2	2.22	0.68
22:CV:38:U:H2'	22:CV:39:C:H6	1.56	0.68
28:D2:26:ARG:HH22	54:DX:6:ASP:CA	2.04	0.68
36:DB:24:G:H4'	36:DB:25:A:C8	2.28	0.68
38:DD:144:ALA:HB3	38:DD:192:THR:CG2	2.22	0.68
35:DA:2820:A:H62	39:DE:192:ASN:HB2	1.56	0.68
41:DG:125:PHE:CD2	41:DG:131:TYR:HB3	2.28	0.68
41:DG:130:ASN:ND2	41:DG:161:THR:N	2.40	0.68
35:DA:196:A:H5''	46:DP:46:LYS:NZ	2.07	0.68
52:DV:1:MET:HE3	52:DV:45:THR:H	1.57	0.68
56:DZ:98:MET:O	56:DZ:126:VAL:HG22	1.94	0.68
1:AA:1307:U:H2'	1:AA:1308:U:C6	2.29	0.68
1:AA:425:G:H2'	1:AA:426:G:H8	1.58	0.68
1:AA:671:G:O2'	1:AA:672:U:H5'	1.94	0.68
3:AC:16:ARG:NH1	3:AC:16:ARG:HB2	2.08	0.68
4:AD:194:LEU:N	4:AD:194:LEU:HD22	2.08	0.68
10:AJ:3:LYS:HB2	10:AJ:77:PRO:HD3	1.75	0.68
34:B8:8:LYS:O	34:B8:12:LYS:HG3	1.92	0.68
35:BA:1879:C:C3'	35:BA:1880:C:H5''	2.23	0.68
35:BA:1947:C:H2'	35:BA:1948:G:C8	2.28	0.68
35:BA:2304:G:H5'	35:BA:2305:A:OP2	1.93	0.68
35:BA:2543:G:H2'	35:BA:2544:G:C8	2.28	0.68
35:BA:2672:G:C2'	35:BA:2673:G:H5''	2.23	0.68
35:BA:690:G:H2'	35:BA:691:C:H6	1.59	0.68
35:BA:795:C:H2'	35:BA:796:C:C6	2.28	0.68
35:BA:877:U:C2'	35:BA:878:A:H5''	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:133:LEU:HA	38:BD:136:ILE:CD1	2.23	0.68
41:BG:41:GLN:O	41:BG:43:LEU:HD22	1.93	0.68
42:BH:126:PRO:O	42:BH:127:GLU:HG2	1.92	0.68
50:BT:25:GLY:HA2	50:BT:92:GLY:N	2.09	0.68
1:CA:449:C:O2	16:CP:42:ARG:HD2	1.93	0.68
2:CB:162:ILE:HD12	2:CB:162:ILE:O	1.93	0.68
2:CB:39:ILE:HG22	2:CB:40:HIS:O	1.93	0.68
3:CC:73:PRO:HA	3:CC:76:VAL:HG13	1.76	0.68
7:CG:65:ALA:O	7:CG:69:VAL:HG23	1.93	0.68
9:CI:114:TYR:HD2	9:CI:114:TYR:N	1.90	0.68
10:CJ:48:THR:HG22	10:CJ:49:VAL:H	1.58	0.68
11:CK:61:ALA:HB3	11:CK:90:GLY:HA3	1.73	0.68
25:CY:28:LEU:O	25:CY:30:THR:N	2.25	0.68
35:DA:1894:C:O2'	35:DA:1895:C:H5'	1.94	0.68
35:DA:2051:A:H4'	39:DE:141:ILE:CD1	2.23	0.68
26:D0:36:ILE:HD11	35:DA:2355:C:H4'	1.75	0.68
35:DA:571:A:H5'	35:DA:2030:A:N6	2.07	0.68
38:DD:239:ARG:HH21	38:DD:239:ARG:HG3	1.58	0.68
41:DG:45:GLU:C	41:DG:47:LYS:H	1.96	0.68
46:DP:126:VAL:HA	46:DP:145:PRO:HB2	1.74	0.68
46:DP:38:GLN:HG3	46:DP:39:LYS:N	2.05	0.68
50:DT:28:VAL:CG2	50:DT:47:GLY:N	2.57	0.68
44:DN:42:TRP:N	51:DU:64:ARG:NH1	2.41	0.68
52:DV:4:ILE:HD12	52:DV:40:LEU:HD21	1.76	0.68
1:AA:899:C:O5'	1:AA:899:C:H6	1.77	0.68
3:AC:104:GLN:CD	3:AC:105:GLU:H	1.96	0.68
25:AY:130:ARG:HG2	35:BA:1942:C:O2	1.93	0.68
35:BA:1635:G:C8	35:BA:1635:G:H5'	2.28	0.68
35:BA:2092:U:H5	35:BA:2226:C:OP2	1.76	0.68
35:BA:2817:G:H21	35:BA:2836:U:C1'	2.07	0.68
38:BD:68:LYS:HG3	38:BD:68:LYS:O	1.93	0.68
40:BF:150:GLY:HA2	40:BF:172:TRP:CE3	2.28	0.68
43:BI:98:ALA:CB	43:BI:109:ILE:HB	2.23	0.68
43:BI:140:LEU:HD12	43:BI:141:LYS:H	1.59	0.68
56:BZ:151:HIS:HA	56:BZ:170:THR:HA	1.75	0.68
1:CA:192:U:C4'	20:CT:103:GLY:HA2	2.24	0.68
1:CA:693:G:H2'	1:CA:694:A:C8	2.28	0.68
4:CD:10:ARG:O	4:CD:13:ARG:HB3	1.93	0.68
6:CF:67:MET:HB2	6:CF:68:PRO:HD2	1.75	0.68
10:CJ:32:ALA:HB1	10:CJ:75:ILE:HD11	1.74	0.68
1:CA:522:C:N4	12:CL:53:ARG:HH21	1.89	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:66:VAL:HG23	12:CL:67:THR:O	1.94	0.68
19:CS:33:THR:HG23	19:CS:51:VAL:HA	1.76	0.68
25:CY:109:GLU:O	25:CY:112:LYS:HB3	1.94	0.68
33:D7:30:VAL:HG23	33:D7:31:LEU:H	1.57	0.68
33:D7:7:PRO:HB2	35:DA:1309:G:H4'	1.75	0.68
35:DA:1434:A:H2'	35:DA:1435:G:C8	2.28	0.68
35:DA:1600:C:O2'	35:DA:1601:G:H5'	1.93	0.68
35:DA:1719:G:O2'	35:DA:1720:U:H5'	1.93	0.68
35:DA:154(A):C:H5	35:DA:171:G:N1	1.92	0.68
35:DA:2208:A:H1'	35:DA:2219:G:C4	2.29	0.68
35:DA:556:G:H2'	35:DA:557:U:C6	2.28	0.68
35:DA:803:U:O2'	35:DA:804:A:H5'	1.92	0.68
35:DA:943:U:OP2	46:DP:38:GLN:CD	2.31	0.68
39:DE:116:VAL:HG21	39:DE:122:PHE:CG	2.28	0.68
42:DH:88:LEU:HD21	42:DH:165:ALA:HA	1.75	0.68
43:DI:79:ILE:HB	43:DI:81:VAL:HG23	1.75	0.68
50:DT:96:ARG:HG3	50:DT:98:LYS:O	1.93	0.68
54:DX:36:LYS:NZ	54:DX:39:ILE:HA	2.08	0.68
35:DA:58:G:OP1	54:DX:72:LYS:HB3	1.93	0.68
1:AA:974:A:H1'	14:AN:31:ARG:HH21	1.57	0.68
3:AC:6:HIS:NE2	3:AC:8:ILE:HB	2.08	0.68
1:AA:963:G:N2	10:AJ:55:LYS:HZ3	1.91	0.68
10:AJ:8:LEU:HA	10:AJ:96:ILE:HG22	1.76	0.68
12:AL:76:ASN:OD1	12:AL:108:ALA:HB3	1.92	0.68
17:AQ:68:ARG:N	17:AQ:70:ARG:NH1	2.41	0.68
20:AT:15:ARG:O	20:AT:18:GLN:HB2	1.93	0.68
27:B1:48:LYS:O	27:B1:49:VAL:HG23	1.92	0.68
27:B1:86:SER:C	27:B1:89:GLU:CG	2.61	0.68
35:BA:2087:G:O2'	35:BA:2088:G:H5'	1.93	0.68
35:BA:2590:A:O2'	35:BA:2591:C:H5'	1.94	0.68
35:BA:431:U:O2'	35:BA:432:A:H5'	1.93	0.68
35:BA:926:A:H2'	35:BA:927:G:H8	1.57	0.68
38:BD:236:GLY:O	38:BD:237:GLU:HG2	1.94	0.68
41:BG:5:VAL:HG12	41:BG:6:ALA:N	2.05	0.68
43:BI:113:ARG:HA	43:BI:131:LYS:HE3	1.75	0.68
43:BI:145:VAL:HG12	43:BI:146:ALA:N	2.07	0.68
44:BN:32:THR:CG2	44:BN:37:LYS:HB3	2.23	0.68
47:BQ:66:ILE:HG22	47:BQ:104:PHE:CE2	2.28	0.68
50:BT:50:ILE:CG2	50:BT:99:LEU:HD12	2.19	0.68
53:BW:75:TYR:HD1	53:BW:75:TYR:N	1.92	0.68
56:BZ:155:LEU:HD23	56:BZ:156:LYS:H	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1149:C:H2'	1:CA:1150:U:H6	1.59	0.68
1:CA:349:A:O2'	1:CA:350:G:H5'	1.94	0.68
1:CA:930:C:O2'	1:CA:931:C:H5'	1.94	0.68
10:CJ:3:LYS:HB2	10:CJ:77:PRO:HD3	1.74	0.68
12:CL:85:ILE:HD11	12:CL:98:TYR:HB3	1.74	0.68
28:D2:30:ARG:H	28:D2:30:ARG:CD	2.06	0.68
35:DA:571:A:H5'	35:DA:2030:A:H62	1.55	0.68
35:DA:541:C:H2'	35:DA:542:C:C5	2.28	0.68
39:DE:52:LEU:HD23	39:DE:75:VAL:HG23	1.75	0.68
40:DF:46:ARG:HG3	40:DF:48:THR:HG23	1.76	0.68
41:DG:46:ALA:CB	41:DG:88:ILE:HB	2.23	0.68
42:DH:162:ILE:HD12	42:DH:162:ILE:C	2.14	0.68
47:DQ:20:ALA:HB2	47:DQ:99:PRO:CG	2.23	0.68
50:DT:30:VAL:HG12	50:DT:44:ASP:CA	2.24	0.68
56:DZ:144:LEU:HD21	56:DZ:150:LEU:HD11	1.73	0.68
1:AA:1126:U:H2'	1:AA:1127:G:C8	2.28	0.68
1:AA:1292:U:H2'	1:AA:1293:G:C8	2.28	0.68
1:AA:477:A:O2'	1:AA:479:C:H5'	1.93	0.68
3:AC:58:GLU:O	3:AC:64:VAL:HA	1.93	0.68
5:AE:102:ALA:H	5:AE:107:ARG:HH21	1.40	0.68
1:AA:1377:A:H2'	7:AG:7:ALA:HB3	1.76	0.68
8:AH:121:ASP:O	8:AH:125:ARG:HB2	1.94	0.68
33:B7:48:LYS:HD3	33:B7:48:LYS:N	2.07	0.68
35:BA:445:C:H5"	51:BU:3:ARG:HB2	1.74	0.68
28:B2:44:LEU:HD23	35:BA:61:G:H5'	1.76	0.68
46:BP:59:LEU:HA	46:BP:61:ARG:NE	2.08	0.68
34:B8:30:ARG:NH2	46:BP:62:LEU:HD23	2.08	0.68
48:BR:9:LYS:O	48:BR:10:LEU:CG	2.41	0.68
49:BS:17:ARG:HG3	49:BS:18:ILE:CD1	2.24	0.68
1:CA:1291:G:H4'	9:CI:39:GLY:HA3	1.74	0.68
1:CA:498:U:H2'	1:CA:499:A:H5'	1.76	0.68
1:CA:677:U:H3	1:CA:713:G:H22	1.42	0.68
2:CB:54:THR:HG21	2:CB:201:ILE:HD11	1.76	0.68
3:CC:104:GLN:CD	3:CC:105:GLU:H	1.97	0.68
4:CD:5:ILE:HG22	4:CD:6:GLY:H	1.56	0.68
12:CL:8:ASN:ND2	17:CQ:34:LYS:HE2	2.08	0.68
1:CA:235:C:H1'	17:CQ:61:GLU:OE1	1.93	0.68
31:D5:49:CYS:O	31:D5:56:LYS:HB2	1.94	0.68
35:DA:2514:U:H2'	35:DA:2515:C:C6	2.29	0.68
35:DA:2777:G:C5'	35:DA:2778:A:H5'	2.23	0.68
39:DE:60:ASN:OD1	39:DE:62:PRO:HD2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DF:6:VAL:HG23	40:DF:125:LEU:H	1.59	0.68
41:DG:69:ALA:HB3	41:DG:91:ARG:O	1.93	0.68
44:DN:108:PRO:C	44:DN:109:LYS:HD3	2.14	0.68
45:DO:4:PRO:O	45:DO:5:GLN:CB	2.41	0.68
47:DQ:23:GLY:HA2	47:DQ:101:ARG:HB2	1.74	0.68
47:DQ:8:LYS:HG3	47:DQ:9:TYR:N	2.08	0.68
48:DR:82:GLU:O	48:DR:85:PRO:HD2	1.94	0.68
49:DS:26:LEU:HA	49:DS:39:ILE:HG13	1.73	0.68
50:DT:52:ILE:HG22	50:DT:61:PHE:HB2	1.76	0.68
53:DW:64:MET:O	53:DW:65:LEU:HB3	1.94	0.68
54:DX:59:VAL:HG23	54:DX:74:PRO:HD2	1.75	0.68
54:DX:77:LYS:HA	54:DX:77:LYS:HE3	1.76	0.68
1:AA:533:A:O2'	1:AA:534:U:H5''	1.93	0.68
5:AE:131:ILE:H	5:AE:131:ILE:HD13	1.59	0.68
6:AF:21:LEU:O	6:AF:25:ILE:HG12	1.93	0.68
10:AJ:6:ILE:HG22	10:AJ:98:ILE:HG13	1.76	0.68
28:B2:28:LYS:O	28:B2:32:LEU:HD23	1.94	0.68
34:B8:6:THR:HG21	34:B8:63:PRO:HD3	1.75	0.68
35:BA:2739:U:O2'	35:BA:2740:A:H5'	1.93	0.68
38:BD:63:ARG:NH1	38:BD:86:PRO:HD2	2.08	0.68
46:BP:17:LYS:C	46:BP:19:VAL:H	1.96	0.68
46:BP:58:THR:C	46:BP:60:MET:H	1.95	0.68
47:BQ:131:ILE:HD13	47:BQ:131:ILE:N	2.08	0.68
50:BT:32:TYR:O	50:BT:33:LYS:HB2	1.94	0.68
35:BA:1614:A:N6	53:BW:93:ALA:HB2	2.09	0.68
1:CA:1046:A:H3'	1:CA:1047:G:H8	1.59	0.68
1:CA:1260:C:OP1	1:CA:1284:C:H4'	1.94	0.68
1:CA:1287:A:H2'	1:CA:1288:A:C8	2.28	0.68
2:CB:19:HIS:O	2:CB:39:ILE:HG23	1.94	0.68
3:CC:111:LEU:HD21	3:CC:146:ALA:HB2	1.73	0.68
9:CI:65:VAL:HG22	9:CI:66:ARG:N	2.08	0.68
12:CL:85:ILE:HD12	12:CL:99:HIS:O	1.94	0.68
20:CT:15:ARG:O	20:CT:18:GLN:HB2	1.94	0.68
20:CT:63:ILE:HD12	20:CT:81:LYS:HG2	1.74	0.68
34:D8:52:LYS:N	34:D8:53:PRO:CD	2.57	0.68
35:DA:110:G:O2'	35:DA:111:A:H5'	1.92	0.68
35:DA:1643:G:H2'	35:DA:1644:C:H6	1.59	0.68
38:DD:4:LYS:HZ1	38:DD:20:ASP:HA	1.59	0.68
47:DQ:34:LEU:HD12	47:DQ:35:VAL:N	2.08	0.68
56:DZ:3:TYR:O	56:DZ:57:ILE:HA	1.94	0.68
1:AA:1260:C:OP1	1:AA:1284:C:H4'	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1424:C:H2'	1:AA:1425:U:H6	1.58	0.68
1:AA:1458:G:H2'	1:AA:1459:C:H6	1.59	0.68
1:AA:627:G:H2'	1:AA:628:G:H8	1.57	0.68
9:AI:10:ARG:HG3	9:AI:104:ARG:O	1.92	0.68
13:AM:66:LEU:CA	13:AM:70:LEU:HD12	2.23	0.68
25:AY:130:ARG:HH11	25:AY:130:ARG:HG3	1.58	0.68
26:B0:27:GLU:HB2	26:B0:69:PHE:CD1	2.28	0.68
27:B1:58:ILE:CD1	27:B1:59:THR:H	1.95	0.68
34:B8:30:ARG:NH2	46:BP:62:LEU:HB2	2.06	0.68
1:AA:1442(A):G:O6	35:BA:2863:C:H4'	1.94	0.68
39:BE:59:VAL:CG2	39:BE:63:LEU:HA	2.23	0.68
44:BN:30:ILE:HD13	44:BN:54:VAL:HG21	1.74	0.68
46:BP:38:GLN:CG	46:BP:39:LYS:H	2.00	0.68
54:BX:12:VAL:HG12	54:BX:27:THR:O	1.93	0.68
54:BX:77:LYS:HE3	54:BX:77:LYS:HA	1.76	0.68
1:CA:1084:G:H5'	1:CA:1102:A:OP2	1.94	0.68
2:CB:204:ASN:HD21	2:CB:207:ALA:H	1.39	0.68
15:CO:61:GLY:O	15:CO:65:ARG:HD3	1.93	0.68
6:CF:91:VAL:HG11	18:CR:72:ARG:HH12	1.58	0.68
27:D1:13:ILE:HG23	27:D1:14:VAL:N	2.07	0.68
27:D1:60:PHE:CD1	27:D1:70:VAL:HG13	2.26	0.68
27:D1:86:SER:N	27:D1:87:PRO:HD3	2.09	0.68
28:D2:50:ILE:O	28:D2:51:ARG:HB2	1.93	0.68
28:D2:53:LEU:HD23	28:D2:54:LYS:HG3	1.75	0.68
29:D3:7:LYS:O	29:D3:54:VAL:HG13	1.93	0.68
42:DH:156:ALA:C	42:DH:158:HIS:N	2.46	0.68
42:DH:145:ALA:HB1	42:DH:164:TYR:CE1	2.29	0.68
45:DO:69:ILE:CD1	45:DO:77:ILE:HG23	2.23	0.68
52:DV:75:PHE:CD1	52:DV:87:HIS:HB3	2.29	0.68
54:DX:12:VAL:HG12	54:DX:27:THR:O	1.94	0.68
55:DY:31:LEU:CD1	55:DY:34:LYS:H	2.06	0.68
1:AA:989:C:H42	1:AA:1217:C:H42	1.40	0.68
1:AA:977:A:C2'	1:AA:978:A:H5'	2.23	0.68
10:AJ:62:HIS:H	10:AJ:62:HIS:CD2	2.12	0.68
11:AK:21:ILE:HB	11:AK:84:VAL:HG12	1.74	0.68
25:AY:116:ARG:HG2	25:AY:116:ARG:HH11	1.59	0.68
27:B1:20:ARG:HD2	27:B1:20:ARG:N	2.09	0.68
32:B6:40:CYS:HB2	32:B6:46:HIS:HE1	1.59	0.68
35:BA:1249:U:H5'	35:BA:1249:U:H6	1.59	0.68
35:BA:2175:C:C2'	35:BA:2176:A:H5''	2.23	0.68
35:BA:2853:C:H2'	35:BA:2854:G:C8	2.28	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:64:A:O2'	35:BA:65:C:H5'	1.94	0.68
35:BA:649:G:H2'	35:BA:650:C:C6	2.28	0.68
39:BE:77:ILE:CG2	39:BE:78:LEU:H	2.07	0.68
39:BE:77:ILE:HG23	39:BE:78:LEU:HD23	1.75	0.68
40:BF:5:ALA:O	40:BF:6:VAL:HG13	1.93	0.68
41:BG:82:LEU:HD22	41:BG:86:MET:HB2	1.76	0.68
43:BI:88:ILE:HD11	43:BI:123:LEU:HD12	1.76	0.68
43:BI:79:ILE:HB	43:BI:81:VAL:HG23	1.76	0.68
44:BN:3:THR:HG22	44:BN:5:VAL:HG23	1.76	0.68
54:BX:55:ASN:HD22	54:BX:55:ASN:N	1.92	0.68
56:BZ:125:LEU:HB3	56:BZ:165:VAL:HG22	1.74	0.68
1:CA:1249:C:H5''	9:CI:70:LYS:HE2	1.74	0.68
2:CB:212:GLN:NE2	2:CB:216:SER:HB2	2.09	0.68
3:CC:110:ASN:O	3:CC:141:VAL:HG13	1.94	0.68
3:CC:84:ILE:HD11	3:CC:88:ARG:HH21	1.59	0.68
4:CD:142:PRO:HA	4:CD:185:PHE:CD2	2.28	0.68
8:CH:85:ARG:HH12	8:CH:134:ILE:HG23	1.59	0.68
6:CF:62:TRP:HB2	18:CR:35:ARG:HH12	1.58	0.68
27:D1:41:ARG:NH2	35:DA:205:G:H1	1.92	0.68
31:D5:40:LYS:HZ3	31:D5:45:VAL:HA	1.56	0.68
35:DA:1007:C:HO2'	44:DN:108:PRO:HA	1.59	0.68
35:DA:1613:G:H2'	35:DA:1617:C:N4	2.09	0.68
35:DA:189:G:H2'	35:DA:205:G:H22	1.59	0.68
35:DA:2505:G:H2'	35:DA:2576:G:O6	1.94	0.68
35:DA:2853:C:H2'	35:DA:2854:G:C8	2.28	0.68
35:DA:708:C:N4	35:DA:723:G:H1	1.92	0.68
35:DA:860:U:C5	35:DA:917:A:N7	2.62	0.68
35:DA:967:C:O2'	35:DA:968:G:H5'	1.94	0.68
36:DB:79:C:O2'	36:DB:80:U:H5'	1.92	0.68
38:DD:145:VAL:HG12	38:DD:146:GLU:N	2.08	0.68
38:DD:215:LEU:O	38:DD:217:ARG:N	2.27	0.68
43:DI:91:SER:CB	43:DI:121:LYS:HE3	2.21	0.68
43:DI:71:ILE:HG13	43:DI:72:LEU:N	2.09	0.68
46:DP:7:ARG:O	46:DP:10:PRO:HD3	1.94	0.68
48:DR:41:ALA:HB1	48:DR:114:VAL:CG2	2.23	0.68
8:AH:45:ILE:HA	8:AH:64:LYS:HB3	1.76	0.68
1:AA:972:C:H4'	10:AJ:57:LYS:HG3	1.76	0.68
25:AY:30:THR:C	25:AY:32:ARG:H	1.97	0.68
35:BA:2544:G:H2'	35:BA:2545:G:H8	1.59	0.68
35:BA:2814:C:H2'	35:BA:2815:C:H6	1.58	0.68
35:BA:909:A:H1'	47:BQ:10:ARG:NH2	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:979:G:H3'	35:BA:980:A:H5''	1.76	0.68
38:BD:45:ASN:ND2	38:BD:46:GLN:H	1.92	0.68
41:BG:153:ARG:NH1	41:BG:153:ARG:HG3	2.01	0.68
46:BP:90:ARG:HD2	46:BP:91:PHE:CD1	2.29	0.68
47:BQ:16:ARG:C	47:BQ:17:LEU:HD23	2.14	0.68
47:BQ:20:ALA:HB2	47:BQ:99:PRO:CG	2.24	0.68
52:BV:19:LYS:HE2	52:BV:19:LYS:HA	1.75	0.68
54:BX:60:ARG:HG3	54:BX:72:LYS:N	2.07	0.68
47:BQ:63:LYS:HA	56:BZ:178:GLU:OE1	1.93	0.68
1:CA:1030:C:H2'	1:CA:1030(A):G:H5'	1.76	0.68
1:CA:624:C:H4'	16:CP:10:GLY:HA2	1.76	0.68
1:CA:741:G:H2'	1:CA:742:G:C8	2.28	0.68
2:CB:102:LEU:HD12	2:CB:102:LEU:N	2.09	0.68
4:CD:189:PRO:HB2	4:CD:194:LEU:HD21	1.75	0.68
10:CJ:27:ALA:HB2	10:CJ:85:LEU:HD11	1.74	0.68
11:CK:103:LEU:H	11:CK:103:LEU:CD2	2.05	0.68
3:AC:79:ARG:NH1	11:CK:99:GLN:HB3	2.09	0.68
16:CP:38:TYR:O	16:CP:49:LEU:HA	1.94	0.68
17:CQ:86:GLU:C	17:CQ:88:TYR:H	1.97	0.68
23:CW:7:G:H2'	23:CW:50:G:OP2	1.94	0.68
25:CY:38:LEU:HA	25:CY:41:LEU:CD1	2.23	0.68
32:D6:10:LEU:CD1	34:D8:36:LYS:HD3	2.13	0.68
35:DA:2316:C:H1'	41:DG:128:ARG:HH12	1.58	0.68
35:DA:2728:U:O2'	35:DA:2729:G:H5'	1.94	0.68
35:DA:673:C:H5'	40:DF:54:ARG:HH12	1.59	0.68
36:DB:42:C:C6	41:DG:69:ALA:HB2	2.29	0.68
38:DD:27:THR:C	38:DD:29:PRO:HD2	2.14	0.68
38:DD:92:ILE:HA	38:DD:107:ALA:H	1.59	0.68
39:DE:103:ASP:OD2	39:DE:201:THR:HA	1.94	0.68
41:DG:2:PRO:HD2	41:DG:4:ASP:O	1.93	0.68
41:DG:51:ARG:HH11	41:DG:53:LEU:HG	1.54	0.68
41:DG:60:LEU:HA	41:DG:63:ILE:CG1	2.24	0.68
42:DH:149:ARG:HA	42:DH:162:ILE:HD11	1.76	0.68
43:DI:23:PRO:HB3	43:DI:27:ARG:HH21	1.58	0.68
48:DR:12:ARG:HH11	48:DR:12:ARG:HG3	1.59	0.68
50:DT:27:THR:O	50:DT:28:VAL:HB	1.94	0.68
1:AA:1190:G:OP1	3:AC:5:ILE:HD12	1.94	0.67
1:AA:1324:A:H2'	1:AA:1325:C:C6	2.19	0.67
1:AA:284:G:H2'	1:AA:285:G:H8	1.58	0.67
1:AA:679:C:O2'	1:AA:680:C:H5'	1.93	0.67
2:AB:104:ASN:OD1	2:AB:107:THR:HB	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:58:LEU:HD13	4:AD:59:ARG:N	2.10	0.67
10:AJ:51:ARG:H	10:AJ:60:ARG:HA	1.57	0.67
11:AK:59:TYR:O	11:AK:63:LEU:HG	1.94	0.67
13:AM:91:ARG:CB	13:AM:98:VAL:HG22	2.24	0.67
19:AS:62:ILE:HD12	19:AS:63:THR:H	1.58	0.67
20:AT:57:ARG:HB2	20:AT:57:ARG:NH1	2.09	0.67
35:BA:1188:U:C2'	35:BA:1189:A:H5'	2.24	0.67
35:BA:1719:G:O2'	35:BA:1720:U:H5'	1.94	0.67
38:BD:242:ARG:HG3	38:BD:242:ARG:HH11	1.58	0.67
41:BG:17:PRO:HA	41:BG:20:ILE:HG12	1.75	0.67
42:BH:116:GLU:HG2	42:BH:117:PRO:HD2	1.77	0.67
43:BI:129:THR:HA	43:BI:137:PRO:HA	1.75	0.67
35:BA:2684:U:P	50:BT:53:ARG:HH21	2.17	0.67
50:BT:60:THR:HG22	50:BT:77:PRO:HA	1.75	0.67
56:BZ:108:PRO:HB3	56:BZ:142:SER:O	1.93	0.67
1:CA:1091:U:H2'	1:CA:1093:A:OP2	1.93	0.67
1:CA:112:G:H4'	1:CA:389:A:H5''	1.74	0.67
1:CA:1253:G:H2'	1:CA:1254:C:C6	2.29	0.67
1:CA:34:C:H2'	1:CA:35:G:C8	2.28	0.67
1:CA:59:A:H5'	1:CA:60:A:H5''	1.76	0.67
1:CA:724:G:O2'	1:CA:725:G:H5'	1.93	0.67
9:CI:5:TYR:HE1	9:CI:7:THR:HG1	1.42	0.67
14:CN:27:CYS:HB3	14:CN:43:CYS:SG	2.33	0.67
23:CW:6:G:H1	23:CW:68:C:H42	1.42	0.67
25:CY:110:ARG:O	25:CY:114:LEU:HD23	1.93	0.67
31:D5:45:VAL:HG22	31:D5:51:TYR:CD1	2.29	0.67
34:D8:30:ARG:NH2	46:DP:62:LEU:HB2	2.09	0.67
35:DA:1414:G:H2'	35:DA:1415:U:H6	1.59	0.67
35:DA:1456:G:H2'	35:DA:1457:A:C8	2.29	0.67
35:DA:186:G:O2'	35:DA:187:G:H5'	1.94	0.67
35:DA:2777:G:C4'	35:DA:2778:A:H5'	2.25	0.67
38:DD:101:GLU:OE1	38:DD:103:ARG:HD3	1.93	0.67
38:DD:24:ILE:O	38:DD:24:ILE:HG23	1.93	0.67
41:DG:39:ILE:CG1	41:DG:157:ILE:HG22	2.24	0.67
41:DG:92:VAL:HG22	41:DG:93:THR:N	2.08	0.67
43:DI:145:VAL:HG12	43:DI:146:ALA:N	2.09	0.67
47:DQ:28:ALA:CB	47:DQ:67:ARG:HD2	2.23	0.67
1:AA:1202:G:H2'	1:AA:1203:C:H5'	1.76	0.67
2:AB:101:MET:HB2	2:AB:102:LEU:HD12	1.74	0.67
2:AB:102:LEU:CD1	2:AB:102:LEU:H	2.07	0.67
1:AA:1096:C:H5''	2:AB:137:ARG:HH22	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:59:ARG:O	4:AD:62:GLN:HB2	1.93	0.67
6:AF:72:VAL:HG13	6:AF:73:ASN:N	2.08	0.67
7:AG:113:GLU:CG	7:AG:119:ARG:HG2	2.24	0.67
1:AA:600:C:OP1	8:AH:97:VAL:HG12	1.93	0.67
18:AR:50:ILE:CD1	18:AR:70:ILE:HG21	2.23	0.67
26:B0:36:ILE:HD11	35:BA:2355:C:C4'	2.22	0.67
27:B1:71:TYR:HA	27:B1:74:VAL:CG2	2.23	0.67
35:BA:1493:C:H4'	35:BA:1494:A:OP1	1.94	0.67
35:BA:2722:G:H2'	35:BA:2723:C:C6	2.29	0.67
39:BE:167:VAL:HG22	39:BE:168:MET:N	2.09	0.67
41:BG:38:VAL:O	41:BG:158:ALA:HB3	1.94	0.67
43:BI:57:ARG:NH1	43:BI:57:ARG:HB3	2.09	0.67
35:BA:943:U:OP2	46:BP:38:GLN:CD	2.32	0.67
47:BQ:23:GLY:HA2	47:BQ:101:ARG:HB2	1.76	0.67
49:BS:26:LEU:C	49:BS:88:ASP:HB3	2.13	0.67
1:CA:233:C:H2'	1:CA:234:C:H6	1.58	0.67
1:CA:1206:G:H4'	3:CC:192:THR:O	1.94	0.67
5:CE:102:ALA:H	5:CE:107:ARG:HH21	1.42	0.67
5:CE:12:LEU:HD11	5:CE:31:LEU:HD13	1.75	0.67
8:CH:20:TYR:CE2	8:CH:75:ARG:HB3	2.29	0.67
22:CV:39:C:O2'	22:CV:40:C:H5'	1.93	0.67
27:D1:62:VAL:HG22	27:D1:63:ALA:N	2.09	0.67
27:D1:88:LYS:O	27:D1:90:ILE:N	2.28	0.67
31:D5:46:CYS:SG	31:D5:47:PRO:HD2	2.35	0.67
34:D8:46:ARG:O	34:D8:47:LYS:HB3	1.93	0.67
35:DA:137:C:H2'	35:DA:139:G:H5'	1.75	0.67
35:DA:1902:C:H4'	38:DD:244:ARG:HB2	1.76	0.67
35:DA:2304:G:H5'	35:DA:2305:A:OP2	1.94	0.67
35:DA:797:C:H2'	35:DA:798:G:H8	1.60	0.67
39:DE:119:ARG:HD2	39:DE:120:TRP:NE1	2.09	0.67
39:DE:11:MET:CB	39:DE:24:THR:HA	2.23	0.67
41:DG:132:ASN:HD21	41:DG:157:ILE:CG1	2.04	0.67
42:DH:97:ARG:O	42:DH:125:VAL:HG11	1.92	0.67
42:DH:13:LYS:C	42:DH:15:VAL:H	1.98	0.67
47:DQ:132:VAL:HG12	47:DQ:133:ARG:H	1.59	0.67
53:DW:26:GLY:HA2	53:DW:71:VAL:O	1.93	0.67
1:AA:1072:G:H2'	1:AA:1073:U:C6	2.29	0.67
1:AA:1422:G:H2'	1:AA:1423:G:C8	2.29	0.67
1:AA:245:C:O2'	1:AA:246:A:H5'	1.94	0.67
1:AA:59:A:H5'	1:AA:60:A:H5''	1.75	0.67
1:AA:689:C:P	11:AK:46:GLY:HA3	2.33	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:696:A:H2'	1:AA:697:U:C6	2.29	0.67
2:AB:187:LEU:C	2:AB:187:LEU:HD22	2.15	0.67
2:AB:187:LEU:HA	2:AB:201:ILE:HB	1.77	0.67
6:AF:68:PRO:HG3	6:AF:71:ARG:NH2	2.08	0.67
11:AK:65:ALA:HB3	11:AK:97:ALA:HB3	1.76	0.67
1:AA:562:C:H1'	12:AL:15:ARG:HB3	1.75	0.67
12:AL:46:LYS:HZ1	12:AL:47:LYS:HB2	1.58	0.67
12:AL:66:VAL:HG23	12:AL:67:THR:O	1.93	0.67
16:AP:71:ARG:O	16:AP:74:LEU:HB2	1.93	0.67
28:B2:30:ARG:HD2	28:B2:30:ARG:N	2.00	0.67
31:B5:40:LYS:HE2	31:B5:46:CYS:CB	2.25	0.67
35:BA:2838:G:O2'	35:BA:2839:G:H5'	1.94	0.67
35:BA:933:A:H2'	35:BA:934:G:O4'	1.94	0.67
36:BB:61:G:H2'	36:BB:62:C:C6	2.29	0.67
39:BE:52:LEU:HD23	39:BE:75:VAL:HG23	1.77	0.67
39:BE:60:ASN:OD1	39:BE:62:PRO:HD2	1.94	0.67
35:BA:323:G:H2'	40:BF:169:ASN:ND2	2.09	0.67
40:BF:32:LEU:C	40:BF:32:LEU:HD23	2.15	0.67
44:BN:30:ILE:O	44:BN:34:LEU:HB2	1.94	0.67
35:BA:389:G:H1	46:BP:71:VAL:CG1	2.07	0.67
48:BR:116:LEU:O	48:BR:117:VAL:HB	1.94	0.67
1:CA:1288:A:H1'	1:CA:1352:C:O2'	1.94	0.67
1:CA:1457:G:H2'	1:CA:1458:G:H8	1.60	0.67
4:CD:120:LEU:HB3	4:CD:125:HIS:HB2	1.75	0.67
11:CK:27:ASN:HA	11:CK:55:LYS:O	1.95	0.67
13:CM:39:ILE:HD12	13:CM:56:LEU:HD23	1.75	0.67
35:DA:104:U:H2'	35:DA:105:C:O4'	1.94	0.67
35:DA:1177:A:H5'	35:DA:1178:C:C6	2.30	0.67
35:DA:796:C:H2'	35:DA:797:C:C6	2.28	0.67
35:DA:833:U:H2'	35:DA:834:C:C6	2.29	0.67
38:DD:133:LEU:HA	38:DD:136:ILE:CD1	2.24	0.67
41:DG:60:LEU:HD13	41:DG:63:ILE:HG12	1.75	0.67
44:DN:30:ILE:HD13	44:DN:54:VAL:HG21	1.75	0.67
46:DP:131:SER:HB2	46:DP:134:ALA:CB	2.24	0.67
46:DP:83:VAL:CG1	46:DP:112:LEU:HD21	2.25	0.67
47:DQ:55:VAL:HG12	47:DQ:64:ILE:CD1	2.20	0.67
1:AA:357:G:OP1	1:AA:367:U:H5''	1.94	0.67
3:AC:138:VAL:HG22	3:AC:151:VAL:HG23	1.76	0.67
6:AF:12:PRO:O	6:AF:14:LEU:N	2.27	0.67
18:AR:19:LYS:O	18:AR:20:ALA:HB2	1.94	0.67
35:BA:2020:A:O2'	35:BA:2021:C:H5'	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BB:24:G:H4'	36:BB:25:A:C8	2.29	0.67
38:BD:10:THR:HG23	38:BD:13:ARG:CB	2.25	0.67
43:BI:4:ILE:HD13	43:BI:47:LEU:HD22	1.74	0.67
44:BN:108:PRO:C	44:BN:109:LYS:HD3	2.15	0.67
35:BA:1131:G:OP1	44:BN:80:GLY:HA2	1.95	0.67
52:BV:39:LEU:HD11	52:BV:53:GLU:N	2.08	0.67
53:BW:26:GLY:O	53:BW:27:LYS:HG3	1.94	0.67
56:BZ:4:ARG:O	56:BZ:5:LEU:HG	1.94	0.67
9:CI:50:LEU:O	9:CI:53:VAL:HG22	1.94	0.67
11:CK:59:TYR:O	11:CK:63:LEU:HG	1.95	0.67
17:CQ:67:LYS:HA	17:CQ:70:ARG:HH12	1.60	0.67
35:DA:1438:U:H2'	35:DA:1439:A:C8	2.30	0.67
35:DA:1908:C:H2'	35:DA:1909:C:H6	1.58	0.67
35:DA:2817:G:H21	35:DA:2836:U:C1'	2.06	0.67
35:DA:649:G:H2'	35:DA:650:C:C6	2.29	0.67
39:DE:173:VAL:HG12	39:DE:174:ASP:N	2.10	0.67
35:DA:1190:G:H4'	46:DP:35:HIS:HB3	1.75	0.67
49:DS:26:LEU:C	49:DS:88:ASP:HB3	2.14	0.67
1:AA:1375:A:H2'	1:AA:1376:U:H6	1.59	0.67
1:AA:1483:A:H2'	1:AA:1484:C:H5'	1.76	0.67
3:AC:73:PRO:HA	3:AC:76:VAL:CG1	2.25	0.67
4:AD:201:GLN:HA	4:AD:204:ILE:HD12	1.76	0.67
5:AE:109:ILE:HG22	5:AE:110:LEU:N	2.08	0.67
8:AH:10:LEU:HD23	8:AH:10:LEU:H	1.60	0.67
25:AY:28:LEU:HB3	25:AY:114:LEU:HD11	1.76	0.67
26:B0:70:GLN:CG	26:B0:71:ASP:H	2.02	0.67
30:B4:45:GLY:C	30:B4:47:GLN:H	1.98	0.67
33:B7:30:VAL:HG23	33:B7:31:LEU:H	1.59	0.67
34:B8:25:MET:HB2	46:BP:62:LEU:HD11	1.77	0.67
35:BA:1345:C:O2'	35:BA:1346:G:H5'	1.94	0.67
35:BA:1438:U:H2'	35:BA:1439:A:C8	2.29	0.67
38:BD:167:GLY:H	38:BD:174:ILE:HB	1.58	0.67
38:BD:24:ILE:HG23	38:BD:24:ILE:O	1.93	0.67
42:BH:89:ILE:HD11	42:BH:129:THR:HB	1.75	0.67
48:BR:82:GLU:C	48:BR:85:PRO:HD2	2.15	0.67
52:BV:75:PHE:CD1	52:BV:87:HIS:HB3	2.30	0.67
54:BX:36:LYS:NZ	54:BX:39:ILE:HA	2.08	0.67
1:CA:1253:G:H5'	10:CJ:44:VAL:HG12	1.76	0.67
1:CA:426:G:H4'	4:CD:41:GLY:O	1.94	0.67
1:CA:439:A:C2	1:CA:441:A:H1'	2.28	0.67
1:CA:818:G:H3'	1:CA:819:A:C5'	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:22:LYS:HZ2	2:CB:22:LYS:HA	1.59	0.67
3:CC:181:ASN:ND2	3:CC:204:LEU:HB2	2.10	0.67
4:CD:128:VAL:HG12	4:CD:129:ASN:N	2.09	0.67
8:CH:82:HIS:HB3	8:CH:138:TRP:CE2	2.28	0.67
25:CY:129:ILE:HA	25:CY:132:ILE:HD12	1.75	0.67
25:CY:32:ARG:NE	25:CY:32:ARG:HA	2.08	0.67
27:D1:58:ILE:HD12	27:D1:59:THR:N	2.04	0.67
27:D1:83:GLU:OE1	27:D1:86:SER:N	2.27	0.67
35:DA:142(A):C:H2'	35:DA:143:G:O4'	1.94	0.67
35:DA:853:G:H2'	35:DA:854:G:H8	1.59	0.67
35:DA:2820:A:H62	39:DE:192:ASN:CB	2.08	0.67
39:DE:77:ILE:CG2	39:DE:78:LEU:H	2.07	0.67
46:DP:146:VAL:HG22	46:DP:147:LEU:N	2.05	0.67
47:DQ:108:GLY:O	47:DQ:109:VAL:HG23	1.94	0.67
1:AA:920:U:H1'	1:AA:1080:A:H2	1.59	0.67
1:AA:342:C:O2'	1:AA:343:U:H5'	1.95	0.67
4:AD:128:VAL:C	4:AD:130:GLY:H	1.96	0.67
13:AM:39:ILE:HD12	13:AM:56:LEU:HD23	1.76	0.67
10:AJ:49:VAL:CG1	14:AN:41:ARG:HB2	2.25	0.67
20:AT:80:ARG:O	20:AT:84:LEU:HB2	1.94	0.67
25:AY:150:SER:OG	25:AY:153:GLU:HG3	1.95	0.67
26:B0:25:ARG:HB2	26:B0:37:LEU:HD23	1.77	0.67
27:B1:83:GLU:HG3	27:B1:86:SER:N	2.08	0.67
35:BA:1108:U:H2'	35:BA:1109:C:H5'	1.75	0.67
35:BA:1175:U:H4'	35:BA:1176:G:H3'	1.77	0.67
35:BA:2521:C:N4	35:BA:2544:G:H1	1.92	0.67
35:BA:2808:U:C2'	35:BA:2809:A:H5'	2.23	0.67
35:BA:554:U:O2'	35:BA:555:U:H5'	1.95	0.67
35:BA:680:G:H2'	35:BA:681:G:H8	1.56	0.67
35:BA:322:A:P	40:BF:169:ASN:HB2	2.34	0.67
41:BG:76:SER:HB3	41:BG:83:ARG:HB3	1.77	0.67
45:BO:4:PRO:O	45:BO:5:GLN:CB	2.42	0.67
52:BV:14:VAL:HG12	52:BV:15:GLU:H	1.60	0.67
1:CA:1090:U:H2'	1:CA:1091:U:C6	2.30	0.67
1:CA:334:C:O2'	1:CA:335:C:H5'	1.94	0.67
11:CK:65:ALA:HB3	11:CK:97:ALA:HB3	1.75	0.67
15:CO:15:PHE:HB2	15:CO:27:VAL:HG22	1.77	0.67
1:CA:333:G:H4'	20:CT:16:HIS:CE1	2.29	0.67
29:D3:36:VAL:O	29:D3:37:LEU:HD23	1.95	0.67
35:DA:1401:G:H2'	35:DA:1402:C:H6	1.59	0.67
35:DA:2028:U:H2'	35:DA:2029:G:C8	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2732:G:C3'	35:DA:2733:A:H5'	2.23	0.67
35:DA:718:A:H3'	35:DA:719:C:C6	2.30	0.67
38:DD:130:ALA:HB1	38:DD:191:ALA:O	1.93	0.67
44:DN:57:ALA:O	44:DN:58:ASP:O	2.13	0.67
44:DN:70:LYS:O	44:DN:71:ILE:HD13	1.95	0.67
47:DQ:43:THR:OG1	47:DQ:46:GLN:HG3	1.93	0.67
51:DU:92:ARG:HD2	52:DV:11:GLN:HG3	1.76	0.67
56:DZ:57:ILE:N	56:DZ:57:ILE:HD12	2.09	0.67
1:AA:1195:C:H5''	1:AA:1196:U:OP2	1.94	0.67
1:AA:15:G:H4'	5:AE:24:ARG:NH2	2.10	0.67
2:AB:39:ILE:HG22	2:AB:40:HIS:O	1.95	0.67
4:AD:33:MET:CE	4:AD:37:PRO:HA	2.25	0.67
7:AG:86:GLN:CG	23:AW:33:C:H5'	2.25	0.67
8:AH:23:SER:HA	8:AH:63:LEU:CD2	2.24	0.67
9:AI:63:ILE:N	9:AI:63:ILE:HD12	2.09	0.67
27:B1:68:PRO:O	27:B1:70:VAL:N	2.28	0.67
31:B5:41:PRO:HG2	31:B5:44:THR:OG1	1.94	0.67
35:BA:2472:G:H5'	35:BA:2473:U:H5''	1.76	0.67
35:BA:577:G:H2'	35:BA:578:A:C8	2.30	0.67
38:BD:158:ALA:O	38:BD:161:THR:HG23	1.94	0.67
38:BD:223:GLY:O	38:BD:225:ALA:N	2.26	0.67
40:BF:134:GLY:H	40:BF:162:LEU:HD11	1.59	0.67
42:BH:88:LEU:HD21	42:BH:165:ALA:HA	1.76	0.67
35:BA:271(P):C:C5'	43:BI:46:ALA:HB2	2.19	0.67
43:BI:71:ILE:HG13	43:BI:72:LEU:N	2.09	0.67
56:BZ:138:GLU:H	56:BZ:138:GLU:CD	1.96	0.67
1:CA:1079:G:H2'	1:CA:1080:A:C8	2.29	0.67
1:CA:989:C:H42	1:CA:1217:C:H42	1.40	0.67
1:CA:1402:C:H2'	1:CA:1403:C:O4'	1.94	0.67
1:CA:294:U:H2'	1:CA:295:C:C6	2.30	0.67
1:CA:340:U:H2'	1:CA:341:C:H6	1.60	0.67
1:CA:490:G:H2'	1:CA:491:G:H8	1.59	0.67
3:CC:127:ARG:HG2	3:CC:127:ARG:HH11	1.57	0.67
6:CF:69:GLU:O	6:CF:72:VAL:HG12	1.95	0.67
20:CT:80:ARG:O	20:CT:84:LEU:HB2	1.94	0.67
26:D0:39:ARG:HH21	35:DA:2355:C:H1'	1.57	0.67
35:DA:108:U:H2'	35:DA:109:G:H8	1.59	0.67
35:DA:1257:C:O2'	40:DF:84:VAL:HG23	1.95	0.67
35:DA:779:U:H2'	35:DA:780:G:C8	2.30	0.67
35:DA:792:G:C5'	35:DA:793:A:H5'	2.25	0.67
37:DC:22:ILE:HG22	37:DC:25:ALA:HB2	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1792:G:P	38:DD:206:LEU:HB2	2.35	0.67
35:DA:2787:C:O2	39:DE:61:ARG:HD3	1.94	0.67
41:DG:140:ILE:HD12	41:DG:141:PHE:N	2.09	0.67
41:DG:51:ARG:CZ	41:DG:51:ARG:HA	2.24	0.67
45:DO:69:ILE:N	45:DO:69:ILE:HD12	2.09	0.67
48:DR:97:VAL:HG22	48:DR:114:VAL:HG22	1.76	0.67
54:DX:53:LYS:NZ	54:DX:55:ASN:HD21	1.92	0.67
1:AA:189(D):C:H1'	1:AA:189(H):G:C2	2.30	0.67
1:AA:331:G:OP1	1:AA:332:G:H5'	1.95	0.67
1:AA:990:C:H2'	1:AA:991:U:C6	2.30	0.67
5:AE:102:ALA:HB1	5:AE:106:PRO:CG	2.25	0.67
12:AL:89:ARG:HH11	12:AL:90:VAL:N	1.93	0.67
15:AO:61:GLY:O	15:AO:65:ARG:HD3	1.95	0.67
23:AW:54:G:O2'	23:AW:55:5MU:H5''	1.95	0.67
35:BA:1232:G:H2'	35:BA:1233:C:H6	1.59	0.67
35:BA:1531:C:H3'	35:BA:1532:C:C4'	2.25	0.67
35:BA:2243:U:H2'	35:BA:2244:U:C6	2.30	0.67
35:BA:272(D):G:H1	35:BA:364:C:N4	1.93	0.67
35:BA:703:U:H2'	35:BA:704:G:H5'	1.76	0.67
38:BD:130:ALA:HB1	38:BD:191:ALA:O	1.95	0.67
35:BA:2830:G:H5'	39:BE:58:ARG:HH12	1.60	0.67
39:BE:65:GLY:O	39:BE:70:ALA:HB2	1.95	0.67
50:BT:83:ILE:HG13	50:BT:84:GLN:N	2.08	0.67
51:BU:83:LEU:CB	51:BU:88:ILE:HG12	2.24	0.67
52:BV:72:VAL:HA	52:BV:88:ARG:NH1	2.09	0.67
54:BX:77:LYS:HE2	54:BX:78:LYS:HG3	1.77	0.67
56:BZ:72:ARG:O	56:BZ:73:GLN:HB2	1.93	0.67
1:CA:284:G:H2'	1:CA:285:G:H8	1.59	0.67
1:CA:332:G:H2'	1:CA:333:G:H8	1.59	0.67
1:CA:899:C:H6	1:CA:899:C:O5'	1.76	0.67
2:CB:162:ILE:C	2:CB:162:ILE:HD12	2.15	0.67
15:CO:54:ARG:HA	15:CO:57:LEU:HD12	1.77	0.67
16:CP:28:ARG:HH11	16:CP:28:ARG:HG2	1.59	0.67
23:CW:57:C:H2'	23:CW:58:A:C8	2.30	0.67
35:DA:795:C:H2'	35:DA:796:C:C6	2.30	0.67
38:DD:242:ARG:HH11	38:DD:242:ARG:HG3	1.60	0.67
38:DD:34:VAL:O	38:DD:34:VAL:HG13	1.94	0.67
39:DE:52:LEU:HD12	39:DE:53:PRO:HD2	1.76	0.67
41:DG:73:ALA:O	41:DG:85:GLY:HA2	1.94	0.67
45:DO:68:GLU:CD	45:DO:78:ARG:HH11	1.97	0.67
46:DP:23:PRO:HD2	46:DP:33:ARG:CZ	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:32:THR:O	46:DP:33:ARG:HB2	1.93	0.67
35:DA:389:G:H1	46:DP:71:VAL:CG1	2.08	0.67
46:DP:85:LEU:CD2	46:DP:85:LEU:H	2.06	0.67
49:DS:17:ARG:HG3	49:DS:18:ILE:CD1	2.25	0.67
1:AA:1499:A:O2'	1:AA:1500:A:H5'	1.94	0.67
1:AA:312:C:H2'	1:AA:313:A:C8	2.30	0.67
4:AD:101:LEU:HD23	4:AD:121:VAL:HG13	1.76	0.67
5:AE:96:PRO:HA	5:AE:117:ASP:OD2	1.95	0.67
5:AE:149:GLU:O	5:AE:153:LYS:HG2	1.94	0.67
15:AO:54:ARG:HA	15:AO:57:LEU:HD12	1.76	0.67
35:BA:1523:U:H2'	35:BA:1524:G:H8	1.60	0.67
35:BA:1742:G:N7	35:BA:1743:C:C4	2.63	0.67
35:BA:2893:G:H5'	35:BA:2894:G:C5'	2.15	0.67
54:BX:38:GLU:HB3	54:BX:41:ASN:HD21	1.59	0.67
1:CA:1064:G:N2	1:CA:1190:G:H2'	2.10	0.67
6:CF:21:LEU:O	6:CF:25:ILE:HG12	1.94	0.67
7:CG:113:GLU:CG	7:CG:119:ARG:HG2	2.25	0.67
10:CJ:6:ILE:HG22	10:CJ:98:ILE:HG13	1.77	0.67
27:D1:11:ARG:HB3	27:D1:12:PRO:HD2	1.77	0.67
31:D5:41:PRO:HG2	31:D5:44:THR:OG1	1.93	0.67
32:D6:40:CYS:HB2	32:D6:46:HIS:HE1	1.58	0.67
35:DA:1531:C:H3'	35:DA:1532:C:C4'	2.25	0.67
35:DA:1707:G:H2'	35:DA:1708:C:C6	2.30	0.67
35:DA:18:C:H2'	35:DA:19:C:C6	2.30	0.67
35:DA:2087:G:O2'	35:DA:2088:G:H5'	1.95	0.67
39:DE:96:PHE:HA	39:DE:100:GLU:OE1	1.94	0.67
41:DG:43:LEU:H	41:DG:43:LEU:CD1	2.08	0.67
50:DT:3:ARG:HB3	50:DT:6:LEU:HB3	1.77	0.67
51:DU:90:VAL:HG22	52:DV:39:LEU:HD12	1.75	0.67
54:DX:38:GLU:HB3	54:DX:41:ASN:HD21	1.60	0.67
1:AA:1079:G:H2'	1:AA:1080:A:C8	2.30	0.67
4:AD:65:ARG:HD2	4:AD:70:ILE:O	1.94	0.67
8:AH:119:LEU:HG	8:AH:124:ALA:HB2	1.75	0.67
8:AH:127:LEU:O	8:AH:127:LEU:HD13	1.95	0.67
14:AN:34:TYR:C	14:AN:36:PHE:H	1.97	0.67
18:AR:67:ALA:HA	18:AR:70:ILE:HD12	1.77	0.67
1:AA:1340:A:O2'	22:AV:31:U:H5'	1.94	0.67
25:AY:112:LYS:O	25:AY:116:ARG:HG3	1.95	0.67
29:B3:7:LYS:O	29:B3:54:VAL:HG13	1.95	0.67
31:B5:45:VAL:HG22	31:B5:51:TYR:CD1	2.29	0.67
35:BA:2222:G:O2'	35:BA:2223:G:H5'	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2820:A:C8	39:BE:191:PRO:HB2	2.29	0.67
44:BN:91:LEU:HA	44:BN:95:PRO:HB3	1.76	0.67
46:BP:146:VAL:HG22	46:BP:147:LEU:N	2.05	0.67
35:BA:631:A:O2'	46:BP:67:MET:HB3	1.95	0.67
47:BQ:65:PHE:O	47:BQ:66:ILE:HG23	1.94	0.67
47:BQ:82:ARG:HG2	47:BQ:82:ARG:HH11	1.59	0.67
50:BT:48:ILE:HD12	50:BT:48:ILE:N	2.10	0.67
2:CB:187:LEU:HA	2:CB:201:ILE:HB	1.77	0.67
2:CB:33:TYR:HD1	2:CB:43:ASP:HA	1.58	0.67
1:CA:1190:G:H3'	3:CC:3:ASN:OD1	1.95	0.67
4:CD:191:ARG:HH12	4:CD:195:ALA:HA	1.59	0.67
6:CF:43:LEU:H	6:CF:43:LEU:HD12	1.60	0.67
8:CH:11:THR:CG2	8:CH:14:ARG:HH12	2.06	0.67
11:CK:88:GLY:O	11:CK:91:ARG:HB2	1.96	0.67
12:CL:46:LYS:HZ1	12:CL:47:LYS:HB2	1.59	0.67
13:CM:75:ALA:O	13:CM:79:LYS:HG3	1.95	0.67
34:D8:8:LYS:HB3	34:D8:12:LYS:HE3	1.77	0.67
35:DA:1108:U:H2'	35:DA:1109:C:H5'	1.75	0.67
35:DA:1935:G:H3'	35:DA:1962:C:H42	1.58	0.67
35:DA:1946:U:O2'	35:DA:1947:C:H5'	1.95	0.67
35:DA:718:A:H3'	35:DA:719:C:H6	1.58	0.67
35:DA:743:G:O2'	35:DA:744:G:H5'	1.95	0.67
35:DA:78:A:H2'	35:DA:79:G:C8	2.29	0.67
42:DH:89:ILE:HD13	42:DH:89:ILE:N	2.10	0.67
44:DN:9:VAL:CG1	44:DN:39:ARG:HH22	2.06	0.67
44:DN:91:LEU:HA	44:DN:95:PRO:HB3	1.77	0.67
46:DP:95:VAL:CG2	46:DP:125:VAL:HB	2.25	0.67
47:DQ:131:ILE:N	47:DQ:131:ILE:HD13	2.10	0.67
48:DR:9:LYS:O	48:DR:10:LEU:CG	2.42	0.67
48:DR:55:ALA:HB2	48:DR:79:LEU:HD12	1.75	0.67
50:DT:38:ASN:HD22	50:DT:40:THR:N	1.88	0.67
1:AA:1046:A:H3'	1:AA:1047:G:H8	1.60	0.66
1:AA:1253:G:H2'	1:AA:1254:C:C6	2.30	0.66
1:AA:160:A:H1'	1:AA:344:A:N7	2.10	0.66
1:AA:59:A:H5'	1:AA:60:A:C5'	2.24	0.66
1:AA:962:C:H2'	1:AA:963:G:H8	1.58	0.66
2:AB:54:THR:HG21	2:AB:201:ILE:HD11	1.75	0.66
16:AP:5:ARG:C	16:AP:6:LEU:HD12	2.16	0.66
23:AW:7:G:H5'	23:AW:8:U:H5	1.59	0.66
25:AY:14:MET:HE1	25:AY:165:THR:HG23	1.77	0.66
27:B1:87:PRO:C	27:B1:89:GLU:N	2.46	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:B8:43:GLN:C	34:B8:44:LYS:HD2	2.16	0.66
35:BA:1713:U:O2'	35:BA:1714:G:H5'	1.95	0.66
35:BA:465:G:H2'	35:BA:466:A:C8	2.29	0.66
35:BA:603:A:H4'	35:BA:604:G:O5'	1.94	0.66
38:BD:144:ALA:HB3	38:BD:192:THR:CG2	2.25	0.66
35:BA:608:A:OP1	40:BF:100:THR:HG21	1.95	0.66
40:BF:28:ILE:HG22	40:BF:112:MET:HB3	1.76	0.66
41:BG:5:VAL:HG11	41:BG:104:GLU:OE2	1.95	0.66
42:BH:13:LYS:C	42:BH:15:VAL:H	1.98	0.66
46:BP:95:VAL:CG2	46:BP:125:VAL:HB	2.25	0.66
48:BR:55:ALA:HB2	48:BR:79:LEU:HD12	1.77	0.66
1:CA:532:A:H2	1:CA:1207:G:O4'	1.78	0.66
1:CA:377:G:OP1	16:CP:3:LYS:HD3	1.95	0.66
1:CA:679:C:O2'	1:CA:680:C:H5'	1.95	0.66
2:CB:104:ASN:OD1	2:CB:107:THR:HB	1.95	0.66
2:CB:194:PRO:O	2:CB:196:LEU:N	2.29	0.66
3:CC:152:ILE:HA	3:CC:166:GLU:O	1.95	0.66
10:CJ:27:ALA:CB	10:CJ:85:LEU:HD11	2.24	0.66
11:CK:44:SER:OG	11:CK:47:VAL:HG23	1.94	0.66
10:CJ:49:VAL:CG1	14:CN:41:ARG:HB2	2.24	0.66
25:CY:43:VAL:O	25:CY:49:HIS:HA	1.95	0.66
27:D1:88:LYS:O	27:D1:92:LYS:N	2.27	0.66
30:D4:45:GLY:C	30:D4:47:GLN:H	1.97	0.66
33:D7:13:ALA:O	33:D7:17:GLY:HA3	1.94	0.66
34:D8:32:LEU:HB3	34:D8:35:GLN:H	1.58	0.66
35:DA:1899:G:N2	35:DA:1902:C:N4	2.43	0.66
35:DA:2443:C:O2'	35:DA:2444:G:H5'	1.95	0.66
35:DA:323:G:C2'	40:DF:169:ASN:HD21	2.08	0.66
35:DA:779:U:H2'	35:DA:780:G:H8	1.60	0.66
38:DD:142:VAL:HG23	38:DD:192:THR:C	2.16	0.66
38:DD:206:LEU:HA	38:DD:211:ARG:HH12	1.58	0.66
39:DE:105:THR:OG1	39:DE:166:THR:HG22	1.94	0.66
39:DE:4:ILE:CG1	39:DE:28:ALA:HB1	2.25	0.66
41:DG:93:THR:O	41:DG:94:LEU:HD23	1.94	0.66
43:DI:38:LEU:HB2	43:DI:40:THR:HG23	1.75	0.66
46:DP:79:ARG:HH21	46:DP:109:GLY:CA	2.06	0.66
46:DP:90:ARG:HD2	46:DP:91:PHE:CD1	2.29	0.66
54:DX:55:ASN:HB2	54:DX:77:LYS:HD2	1.75	0.66
1:AA:1116:C:C3'	1:AA:1117:G:H5''	2.25	0.66
3:AC:83:ARG:O	3:AC:87:LEU:HG	1.95	0.66
4:AD:13:ARG:HD3	4:AD:38:TYR:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:152:ALA:C	7:AG:154:TYR:H	1.97	0.66
14:AN:27:CYS:HB3	14:AN:43:CYS:SG	2.34	0.66
17:AQ:69:LYS:C	17:AQ:70:ARG:HD2	2.15	0.66
25:AY:7:TYR:OH	25:AY:157:ALA:HA	1.94	0.66
35:BA:108:U:H2'	35:BA:109:G:C8	2.30	0.66
35:BA:2845:G:O2'	35:BA:2846:G:H5'	1.95	0.66
40:BF:2:LYS:O	40:BF:25:PRO:HD2	1.95	0.66
43:BI:133:HIS:O	43:BI:135:GLU:HG3	1.94	0.66
47:BQ:30:GLY:CA	47:BQ:107:ALA:HB2	2.24	0.66
47:BQ:108:GLY:O	47:BQ:109:VAL:HG23	1.94	0.66
50:BT:118:ARG:O	50:BT:121:ILE:HG22	1.95	0.66
55:BY:87:LYS:HG3	55:BY:89:PHE:N	2.06	0.66
1:CA:1488:G:H2'	1:CA:1489:G:C8	2.30	0.66
1:CA:59:A:H5'	1:CA:60:A:C5'	2.26	0.66
3:CC:23:TYR:HA	10:CJ:11:PHE:CE1	2.31	0.66
4:CD:22:LYS:HB2	4:CD:26:CYS:CB	2.20	0.66
6:CF:9:VAL:C	6:CF:10:LEU:HD12	2.15	0.66
11:CK:65:ALA:O	11:CK:68:ALA:HB3	1.95	0.66
25:CY:45:TYR:HB2	25:CY:78:ALA:HB1	1.75	0.66
27:D1:37:ILE:HD12	35:DA:2080:G:O5'	1.94	0.66
27:D1:46:LEU:H	27:D1:46:LEU:CD1	1.99	0.66
35:DA:1742:G:N7	35:DA:1743:C:C4	2.63	0.66
35:DA:1778:U:C5	35:DA:1784:A:C4	2.82	0.66
35:DA:2037:G:H2'	35:DA:2038:G:C8	2.30	0.66
35:DA:2115:G:H4'	35:DA:2166:G:N2	2.09	0.66
35:DA:2830:G:H5'	39:DE:58:ARG:HH22	1.61	0.66
35:DA:559:G:N2	51:DU:49:HIS:CD2	2.63	0.66
36:DB:60:C:H2'	36:DB:61:G:H8	1.61	0.66
38:DD:95:LEU:HD12	38:DD:103:ARG:O	1.92	0.66
38:DD:25:THR:CG2	38:DD:82:ILE:H	2.08	0.66
38:DD:25:THR:HG22	38:DD:82:ILE:O	1.95	0.66
45:DO:61:VAL:HG22	45:DO:62:VAL:O	1.95	0.66
47:DQ:10:ARG:HD3	47:DQ:12:GLN:HB3	1.75	0.66
49:DS:62:LYS:HD3	49:DS:62:LYS:N	2.10	0.66
50:DT:28:VAL:CG2	50:DT:47:GLY:H	2.03	0.66
51:DU:105:VAL:O	51:DU:109:LEU:HG	1.95	0.66
55:DY:76:CYS:O	55:DY:78:ALA:N	2.28	0.66
1:AA:484:G:H4'	1:AA:485:G:O5'	1.95	0.66
3:AC:84:ILE:HD11	3:AC:88:ARG:HH21	1.60	0.66
5:AE:144:THR:O	5:AE:148:VAL:HG23	1.96	0.66
8:AH:20:TYR:CE2	8:AH:75:ARG:HB3	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AN:8:GLU:CB	14:AN:12:ARG:HH11	2.09	0.66
17:AQ:86:GLU:C	17:AQ:88:TYR:H	1.96	0.66
22:AV:34:A:H2'	22:AV:35:A:H8	1.58	0.66
23:AW:1:C:O2'	23:AW:2:G:H5'	1.95	0.66
25:AY:101:ILE:H	25:AY:101:ILE:HD12	1.61	0.66
28:B2:14:ARG:NE	28:B2:14:ARG:N	2.43	0.66
35:BA:1456:G:H2'	35:BA:1457:A:C8	2.30	0.66
35:BA:2259:G:H1'	35:BA:2427:C:C2	2.30	0.66
35:BA:2485:G:O2'	35:BA:2486:G:H5'	1.96	0.66
35:BA:271(V):G:H2'	35:BA:271(W):G:O4'	1.95	0.66
35:BA:2774:C:H2'	35:BA:2775:A:C8	2.31	0.66
35:BA:2781:A:H5''	35:BA:2782:G:H5'	1.77	0.66
35:BA:377:C:H2'	35:BA:378:C:H6	1.60	0.66
35:BA:558:G:H2'	35:BA:559:G:H8	1.58	0.66
35:BA:1819:A:OP1	38:BD:161:THR:HG21	1.95	0.66
45:BO:69:ILE:HD12	45:BO:69:ILE:N	2.11	0.66
49:BS:25:ARG:NH2	49:BS:89:ARG:HH12	1.91	0.66
49:BS:54:LEU:HD13	49:BS:58:LEU:O	1.96	0.66
50:BT:53:ARG:HG2	50:BT:53:ARG:HH11	1.58	0.66
52:BV:61:VAL:HG12	52:BV:62:LEU:N	2.11	0.66
54:BX:80:ILE:O	54:BX:81:VAL:HB	1.94	0.66
56:BZ:110:GLY:O	56:BZ:111:VAL:HG12	1.95	0.66
1:CA:1073:U:H2'	1:CA:1074:G:C8	2.29	0.66
1:CA:1405:G:N2	1:CA:1517:G:H22	1.93	0.66
1:CA:425:G:H2'	1:CA:426:G:H8	1.61	0.66
1:CA:671:G:O2'	1:CA:672:U:H5'	1.94	0.66
3:CC:164:ARG:HB2	3:CC:164:ARG:NH1	2.06	0.66
6:CF:18:GLN:O	6:CF:21:LEU:HB2	1.95	0.66
9:CI:77:ILE:O	9:CI:81:ILE:HG12	1.95	0.66
10:CJ:50:ILE:HA	10:CJ:60:ARG:HG2	1.78	0.66
20:CT:41:ILE:C	20:CT:43:LEU:H	1.97	0.66
25:CY:131:ASN:O	25:CY:132:ILE:C	2.33	0.66
33:D7:9:ARG:NH1	35:DA:1309:G:H3'	2.10	0.66
35:DA:2485:G:O2'	35:DA:2486:G:H5'	1.96	0.66
38:DD:25:THR:O	38:DD:26:LYS:HD2	1.95	0.66
38:DD:94:LEU:N	38:DD:94:LEU:HD12	2.06	0.66
40:DF:9:ILE:HG12	40:DF:14:PRO:HA	1.75	0.66
45:DO:1:MET:H3	45:DO:1:MET:HE2	1.60	0.66
50:DT:42:ILE:HG13	50:DT:42:ILE:O	1.94	0.66
54:DX:25:LYS:NZ	54:DX:87:GLN:N	2.38	0.66
1:AA:1203:C:OP1	14:AN:3:ARG:HD2	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:148:G:O2'	1:AA:149:A:H5'	1.95	0.66
1:AA:54:C:H41	1:AA:352:C:H2'	1.60	0.66
1:AA:60:A:P	1:AA:60:A:H8	2.17	0.66
1:AA:639:G:O2'	1:AA:640:A:H5'	1.96	0.66
1:AA:818:G:H3'	1:AA:819:A:C5'	2.25	0.66
1:AA:943:U:H6	1:AA:943:U:O5'	1.79	0.66
3:AC:152:ILE:HA	3:AC:166:GLU:O	1.96	0.66
6:AF:39:LYS:HG2	6:AF:40:VAL:H	1.60	0.66
7:AG:65:ALA:O	7:AG:69:VAL:HG23	1.95	0.66
12:AL:86:ARG:HG2	12:AL:87:GLY:N	2.11	0.66
35:BA:1177:A:H5'	35:BA:1178:C:C6	2.29	0.66
35:BA:1193:G:H2'	35:BA:1194:A:O4'	1.94	0.66
35:BA:1613:G:H2'	35:BA:1617:C:N4	2.10	0.66
31:B5:4:HIS:O	35:BA:2056:G:N2	2.28	0.66
35:BA:2389:G:H5''	35:BA:2390:U:O4'	1.96	0.66
35:BA:2529:G:OP2	35:BA:2530:A:H5''	1.95	0.66
35:BA:2745:C:H2'	35:BA:2746:U:C6	2.30	0.66
35:BA:449:A:O2'	35:BA:450:G:H5'	1.96	0.66
38:BD:268:ARG:HH11	38:BD:268:ARG:HB2	1.61	0.66
38:BD:34:VAL:O	38:BD:34:VAL:HG13	1.95	0.66
54:BX:53:LYS:NZ	54:BX:55:ASN:HD21	1.93	0.66
2:CB:137:ARG:HH11	2:CB:137:ARG:HG2	1.61	0.66
6:CF:37:VAL:HA	6:CF:65:VAL:HG12	1.77	0.66
10:CJ:40:LEU:HB2	10:CJ:41:PRO:HD2	1.76	0.66
13:CM:91:ARG:CB	13:CM:98:VAL:HG22	2.26	0.66
14:CN:4:LYS:O	14:CN:7:ILE:HG12	1.95	0.66
17:CQ:68:ARG:N	17:CQ:70:ARG:NH1	2.43	0.66
25:CY:73:GLN:HB2	25:CY:77:LYS:NZ	2.10	0.66
26:D0:25:ARG:HA	26:D0:29:GLN:NE2	2.10	0.66
27:D1:87:PRO:HB2	27:D1:91:LYS:HZ2	1.61	0.66
35:DA:1547:C:O2'	35:DA:1548:C:H5'	1.95	0.66
35:DA:703:U:C2'	35:DA:704:G:H5'	2.26	0.66
35:DA:892:G:H1	35:DA:894:C:N4	1.93	0.66
42:DH:44:VAL:HG12	42:DH:45:VAL:H	1.61	0.66
42:DH:89:ILE:HD11	42:DH:129:THR:HB	1.76	0.66
47:DQ:30:GLY:CA	47:DQ:107:ALA:HB2	2.24	0.66
48:DR:18:LEU:HD11	48:DR:22:ARG:NH2	2.11	0.66
48:DR:63:ARG:O	48:DR:67:LEU:HD23	1.94	0.66
50:DT:107:ASP:OD1	50:DT:109:GLU:HB2	1.94	0.66
50:DT:27:THR:OG1	50:DT:87:ASP:HA	1.94	0.66
53:DW:75:TYR:N	53:DW:75:TYR:HD1	1.93	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:DX:60:ARG:HG3	54:DX:72:LYS:N	2.09	0.66
56:DZ:58:VAL:HG22	56:DZ:68:PRO:HA	1.77	0.66
56:DZ:61:LEU:HB2	56:DZ:65:GLN:HB2	1.78	0.66
1:AA:1064:G:N2	1:AA:1190:G:H2'	2.10	0.66
1:AA:1434:A:O2'	1:AA:1435:G:H5'	1.94	0.66
1:AA:523:A:N1	12:AL:92:ASP:HB2	2.11	0.66
1:AA:552:U:O2'	1:AA:553:A:H5'	1.94	0.66
2:AB:100:GLY:HA2	2:AB:103:THR:HB	1.78	0.66
3:AC:109:PRO:HA	3:AC:115:LEU:CD1	2.26	0.66
4:AD:12:CYS:SG	4:AD:19:LEU:O	2.54	0.66
5:AE:12:LEU:HD11	5:AE:31:LEU:HD13	1.77	0.66
8:AH:82:HIS:HB3	8:AH:138:TRP:CE2	2.30	0.66
10:AJ:6:ILE:HA	10:AJ:97:GLU:O	1.95	0.66
1:AA:995:C:O2	14:AN:4:LYS:HE2	1.94	0.66
32:B6:13:CYS:HB3	32:B6:49:HIS:HB3	1.76	0.66
35:BA:1665:A:H2'	35:BA:1666:G:O4'	1.95	0.66
35:BA:1902:C:H4'	38:BD:244:ARG:HB2	1.76	0.66
36:BB:7:G:H3'	36:BB:8:U:C5'	2.23	0.66
38:BD:27:THR:C	38:BD:29:PRO:HD2	2.16	0.66
39:BE:49:LEU:HD22	39:BE:49:LEU:N	2.09	0.66
40:BF:41:LEU:HD11	40:BF:184:TYR:HE1	1.61	0.66
40:BF:52:LYS:HD3	40:BF:57:VAL:HA	1.76	0.66
44:BN:70:LYS:O	44:BN:71:ILE:HD13	1.96	0.66
50:BT:100:TYR:HD2	50:BT:103:ARG:CZ	2.08	0.66
53:BW:26:GLY:HA2	53:BW:71:VAL:O	1.95	0.66
56:BZ:39:VAL:HG21	56:BZ:44:PHE:HD2	1.61	0.66
1:CA:817:C:H42	1:CA:1529:G:H1	1.43	0.66
5:CE:78:HIS:CE1	5:CE:143:ARG:H	2.11	0.66
6:CF:72:VAL:HG13	6:CF:73:ASN:N	2.10	0.66
14:CN:8:GLU:CB	14:CN:12:ARG:HH11	2.08	0.66
16:CP:22:THR:HG22	16:CP:32:TYR:HA	1.78	0.66
19:CS:22:LEU:O	19:CS:22:LEU:HD13	1.95	0.66
23:CW:17:C:C4'	23:CW:62:C:H5'	2.26	0.66
25:CY:4:LYS:HA	25:CY:4:LYS:HE3	1.77	0.66
26:D0:36:ILE:HD11	35:DA:2355:C:C4'	2.26	0.66
35:DA:2259:G:H1'	35:DA:2427:C:C2	2.30	0.66
35:DA:2795:G:N3	35:DA:2795:G:H2'	2.10	0.66
40:DF:67:GLN:O	40:DF:68:LYS:HG2	1.94	0.66
45:DO:1:MET:HG3	45:DO:32:TYR:HD2	1.60	0.66
45:DO:68:GLU:HB3	45:DO:78:ARG:HD3	1.78	0.66
50:DT:10:VAL:C	50:DT:12:SER:N	2.45	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:DU:92:ARG:HG3	51:DU:94:ASN:HB3	1.75	0.66
3:AC:173:VAL:O	3:AC:175:LEU:N	2.28	0.66
1:AA:1191:A:H5''	3:AC:4:LYS:HZ3	1.60	0.66
4:AD:134:ASP:O	4:AD:136:PRO:HD3	1.96	0.66
6:AF:43:LEU:H	6:AF:43:LEU:HD12	1.60	0.66
33:B7:7:PRO:HB2	35:BA:1309:G:H4'	1.77	0.66
35:BA:189:G:H2'	35:BA:205:G:N2	2.10	0.66
35:BA:2014:A:H4'	53:BW:94:ASP:OD1	1.95	0.66
35:BA:2186:G:C3'	35:BA:2187:G:H5''	2.26	0.66
35:BA:2723:C:C2'	35:BA:2724:C:H5'	2.26	0.66
35:BA:581:C:H2'	35:BA:582:G:C8	2.30	0.66
35:BA:860:U:C5	35:BA:917:A:N7	2.64	0.66
37:BC:22:ILE:HG22	37:BC:25:ALA:HB2	1.76	0.66
38:BD:209:ALA:O	38:BD:212:SER:HB3	1.94	0.66
40:BF:6:VAL:HG23	40:BF:125:LEU:H	1.59	0.66
1:CA:189(D):C:H1'	1:CA:189(H):G:C2	2.30	0.66
1:CA:243:A:H4'	1:CA:244:U:O5'	1.96	0.66
1:CA:484:G:H4'	1:CA:485:G:O5'	1.95	0.66
3:CC:112:SER:HB3	3:CC:115:LEU:HD12	1.77	0.66
4:CD:17:VAL:HG12	4:CD:18:LYS:N	2.10	0.66
10:CJ:62:HIS:H	10:CJ:62:HIS:CD2	2.11	0.66
28:D2:30:ARG:CZ	28:D2:30:ARG:HB3	2.26	0.66
28:D2:52:ASP:O	28:D2:56:GLN:HG2	1.96	0.66
31:D5:40:LYS:HE2	31:D5:46:CYS:N	2.09	0.66
35:DA:1175:U:H4'	35:DA:1176:G:H3'	1.77	0.66
35:DA:1568:G:H5''	38:DD:61:LEU:HB2	1.78	0.66
35:DA:2101:G:C2	35:DA:2102:U:H1'	2.31	0.66
35:DA:2881:C:H2'	35:DA:2882:A:H8	1.60	0.66
35:DA:543:C:N4	35:DA:551:G:H1	1.93	0.66
35:DA:979:G:H3'	35:DA:980:A:H5''	1.76	0.66
38:DD:131:LEU:CD1	38:DD:136:ILE:HG12	2.26	0.66
38:DD:79:VAL:HG12	38:DD:113:VAL:HA	1.77	0.66
39:DE:133:LYS:C	39:DE:134:ILE:HD13	2.15	0.66
39:DE:65:GLY:O	39:DE:70:ALA:HB2	1.94	0.66
41:DG:44:GLY:C	41:DG:46:ALA:H	1.99	0.66
43:DI:88:ILE:CG2	43:DI:89:TYR:N	2.58	0.66
44:DN:3:THR:HG22	44:DN:5:VAL:CG2	2.25	0.66
50:DT:80:SER:CB	50:DT:81:PRO:CD	2.65	0.66
52:DV:72:VAL:HA	52:DV:88:ARG:NH1	2.10	0.66
53:DW:5:ALA:HB2	53:DW:54:ALA:HB2	1.77	0.66
1:AA:1073:U:H2'	1:AA:1074:G:C8	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:976:G:H22	1:AA:1362:C:H2'	1.59	0.66
1:AA:1392:G:H21	1:AA:1502:A:H8	1.42	0.66
1:AA:1507:A:C2	1:AA:1530:G:C1'	2.79	0.66
1:AA:160:A:H1'	1:AA:344:A:C8	2.31	0.66
1:AA:741:G:H2'	1:AA:742:G:C8	2.30	0.66
1:AA:954:G:H2'	1:AA:955:U:C6	2.30	0.66
4:AD:189:PRO:HB2	4:AD:194:LEU:HD21	1.78	0.66
9:AI:28:VAL:CG1	9:AI:64:THR:HA	2.25	0.66
23:AW:40:C:H2'	23:AW:41:C:C6	2.31	0.66
26:B0:77:ARG:NH2	35:BA:857:C:H5'	2.10	0.66
31:B5:15:ARG:O	31:B5:18:ALA:HB3	1.96	0.66
31:B5:46:CYS:HB3	31:B5:48:GLU:OE2	1.96	0.66
35:BA:1771:C:H2'	35:BA:1772:G:C8	2.30	0.66
35:BA:2570:G:H2'	35:BA:2571:C:C6	2.30	0.66
35:BA:265:A:H1'	35:BA:266:G:O4'	1.95	0.66
35:BA:2752:C:H5	35:BA:2753:A:H62	1.42	0.66
35:BA:2772:C:H2'	35:BA:2773:C:C6	2.31	0.66
35:BA:2777:G:C5'	35:BA:2778:A:H5'	2.25	0.66
38:BD:121:PRO:HB3	38:BD:135:PHE:CE2	2.31	0.66
42:BH:149:ARG:HG3	42:BH:162:ILE:HD11	1.77	0.66
44:BN:26:LEU:CG	44:BN:30:ILE:HD11	2.26	0.66
35:BA:17:G:H4'	51:BU:25:TRP:CZ3	2.30	0.66
53:BW:1:MET:HG3	53:BW:2:GLU:H	1.61	0.66
54:BX:82:GLN:CG	54:BX:83:VAL:H	2.08	0.66
56:BZ:119:GLU:HG3	56:BZ:119:GLU:O	1.96	0.66
56:BZ:17:ALA:O	56:BZ:20:ARG:HB3	1.96	0.66
56:BZ:9:TYR:OH	56:BZ:35:ARG:HG3	1.96	0.66
1:CA:160:A:H1'	1:CA:344:A:C8	2.31	0.66
1:CA:60:A:P	1:CA:60:A:H8	2.18	0.66
1:CA:627:G:H2'	1:CA:628:G:H8	1.59	0.66
1:CA:1096:C:H5''	2:CB:137:ARG:HH22	1.60	0.66
2:CB:25:ASN:OD1	2:CB:27:LYS:HB2	1.95	0.66
7:CG:49:ILE:O	7:CG:49:ILE:HG22	1.95	0.66
8:CH:119:LEU:HG	8:CH:124:ALA:HB2	1.77	0.66
15:CO:30:ALA:O	15:CO:33:THR:HB	1.94	0.66
16:CP:71:ARG:O	16:CP:74:LEU:HB2	1.95	0.66
20:CT:100:ILE:HG22	20:CT:102:GLY:H	1.61	0.66
34:D8:22:VAL:HG21	34:D8:56:GLU:HB2	1.76	0.66
35:DA:1324:G:H3'	35:DA:1325:G:H4'	1.75	0.66
35:DA:2243:U:H2'	35:DA:2244:U:C6	2.30	0.66
35:DA:2562:U:C2'	35:DA:2563:U:H5'	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:268:C:N4	35:DA:424:G:H1	1.94	0.66
35:DA:2715:C:H2'	35:DA:2716:U:H6	1.60	0.66
35:DA:2864:G:H8	35:DA:2864:G:H5'	1.61	0.66
35:DA:493:G:C3'	35:DA:494:G:H5''	2.25	0.66
35:DA:581:C:H2'	35:DA:582:G:C8	2.27	0.66
35:DA:882:G:H22	35:DA:895:U:H3	1.44	0.66
38:DD:142:VAL:HG23	38:DD:193:VAL:CA	2.24	0.66
41:DG:76:SER:HB3	41:DG:83:ARG:CA	2.26	0.66
41:DG:76:SER:CB	41:DG:84:LYS:H	2.08	0.66
46:DP:23:PRO:HD2	46:DP:33:ARG:NH1	2.10	0.66
54:DX:60:ARG:HG2	54:DX:74:PRO:CD	2.25	0.66
56:DZ:30:ASN:C	56:DZ:32:HIS:H	1.99	0.66
1:AA:1504:G:H3'	1:AA:1504:G:P	2.35	0.66
3:AC:181:ASN:ND2	3:AC:204:LEU:HB2	2.11	0.66
4:AD:200:GLU:O	4:AD:204:ILE:HG13	1.95	0.66
6:AF:69:GLU:O	6:AF:72:VAL:HG12	1.95	0.66
8:AH:18:ARG:N	8:AH:78:GLN:NE2	2.44	0.66
9:AI:114:TYR:N	9:AI:114:TYR:HD2	1.92	0.66
10:AJ:5:ARG:HA	10:AJ:73:ASP:OD1	1.95	0.66
23:AW:64:G:H2'	23:AW:65:G:H8	1.60	0.66
25:AY:14:MET:HA	25:AY:132:ILE:HD13	1.76	0.66
35:BA:2079:U:H3	35:BA:2241:A:H61	1.44	0.66
35:BA:2777:G:C4'	35:BA:2778:A:H5'	2.26	0.66
35:BA:543:C:N4	35:BA:551:G:H1	1.94	0.66
35:BA:78:A:H2'	35:BA:79:G:C8	2.31	0.66
46:BP:62:LEU:N	46:BP:62:LEU:HD13	2.11	0.66
49:BS:87:PHE:CG	49:BS:88:ASP:N	2.62	0.66
51:BU:6:THR:O	51:BU:8:VAL:N	2.22	0.66
51:BU:92:ARG:HG3	51:BU:94:ASN:HB3	1.78	0.66
52:BV:19:LYS:HG3	52:BV:20:LEU:N	2.11	0.66
1:CA:1422:G:H4'	45:DO:49:ARG:HH12	1.59	0.66
1:CA:1472:U:O2'	1:CA:1473:A:H5'	1.95	0.66
1:CA:148:G:O2'	1:CA:149:A:H5'	1.95	0.66
1:CA:160:A:H1'	1:CA:344:A:N7	2.10	0.66
2:CB:31:TYR:N	2:CB:31:TYR:CD2	2.62	0.66
3:CC:15:THR:CG2	3:CC:16:ARG:HH12	2.04	0.66
3:CC:83:ARG:O	3:CC:87:LEU:HG	1.96	0.66
6:CF:62:TRP:CZ2	6:CF:64:GLN:HB2	2.30	0.66
8:CH:23:SER:HA	8:CH:63:LEU:CD2	2.21	0.66
16:CP:6:LEU:HB3	16:CP:17:TYR:HD2	1.59	0.66
21:CU:8:THR:O	21:CU:12:LYS:HB2	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:CY:131:ASN:O	25:CY:134:ARG:N	2.28	0.66
27:D1:19:GLN:HE21	35:DA:379:G:H21	1.42	0.66
27:D1:85:LEU:C	27:D1:87:PRO:HD3	2.17	0.66
34:D8:40:GLU:OE1	34:D8:44:LYS:HE3	1.96	0.66
35:DA:703:U:H2'	35:DA:704:G:H5'	1.76	0.66
35:DA:926:A:H2'	35:DA:927:G:H8	1.60	0.66
40:DF:129:PHE:HA	40:DF:142:TRP:NE1	2.10	0.66
40:DF:150:GLY:HA2	40:DF:172:TRP:CE3	2.30	0.66
41:DG:76:SER:O	41:DG:77:ILE:HG23	1.95	0.66
43:DI:28:ASN:O	43:DI:32:PRO:HG2	1.95	0.66
45:DO:104:ARG:HH12	50:DT:35:LYS:HD3	1.59	0.66
49:DS:28:VAL:HG12	49:DS:29:PHE:N	2.05	0.66
52:DV:61:VAL:HG12	52:DV:62:LEU:N	2.10	0.66
2:AB:185:ILE:CG2	2:AB:199:TYR:HB2	2.26	0.66
6:AF:22:GLU:O	6:AF:26:ILE:HG13	1.96	0.66
8:AH:64:LYS:O	8:AH:79:VAL:HG21	1.95	0.66
35:BA:154(A):C:H5	35:BA:171:G:N1	1.94	0.66
35:BA:1991:U:H2'	35:BA:1992:G:C5'	2.26	0.66
35:BA:2101:G:C2	35:BA:2102:U:H1'	2.31	0.66
35:BA:2712:U:H1'	35:BA:2712(A):A:H8	1.61	0.66
35:BA:2866:U:C6	35:BA:2868:A:H1'	2.31	0.66
35:BA:323:G:C2'	40:BF:169:ASN:HD21	2.09	0.66
35:BA:718:A:H3'	35:BA:719:C:H6	1.61	0.66
38:BD:31:LYS:HA	38:BD:31:LYS:HZ1	1.60	0.66
49:BS:31:SER:OG	49:BS:32:LEU:N	2.29	0.66
50:BT:109:GLU:HA	50:BT:112:ARG:HG3	1.77	0.66
56:BZ:97:GLU:O	56:BZ:98:MET:HB3	1.96	0.66
1:CA:1048:G:H4'	14:CN:2:ALA:N	2.10	0.66
1:CA:1152:A:H5'	10:CJ:13:HIS:CD2	2.31	0.66
1:CA:696:A:H2'	1:CA:697:U:H6	1.60	0.66
2:CB:48:MET:HA	2:CB:51:LEU:HD12	1.78	0.66
6:CF:12:PRO:O	6:CF:14:LEU:N	2.28	0.66
7:CG:84:ASN:HD22	7:CG:84:ASN:N	1.92	0.66
9:CI:10:ARG:HG3	9:CI:104:ARG:O	1.96	0.66
12:CL:55:VAL:HG12	12:CL:56:ALA:N	2.06	0.66
15:CO:17:ARG:HG3	15:CO:17:ARG:HH11	1.61	0.66
28:D2:52:ASP:C	28:D2:54:LYS:N	2.48	0.66
35:DA:237:C:H2'	35:DA:238:C:H6	1.61	0.66
35:DA:999:U:O2	35:DA:999:U:H2'	1.95	0.66
44:DN:62:VAL:HG22	44:DN:66:LYS:HG3	1.78	0.66
44:DN:66:LYS:O	44:DN:67:LEU:HD23	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:38:HIS:O	51:DU:67:ALA:HB1	1.96	0.66
1:AA:1152:A:H5''	10:AJ:13:HIS:CD2	2.31	0.66
1:AA:1206:G:O4'	3:AC:194:GLY:HA2	1.96	0.66
1:AA:1433:A:C2	1:AA:1434:A:C4	2.84	0.66
1:AA:1457:G:H2'	1:AA:1458:G:H8	1.60	0.66
1:AA:1479:C:H2'	1:AA:1480:G:H8	1.60	0.66
1:AA:451:A:H1'	1:AA:452:A:C8	2.31	0.66
2:AB:70:PHE:O	2:AB:92:TYR:HB2	1.95	0.66
6:AF:67:MET:HB2	6:AF:68:PRO:HD2	1.77	0.66
8:AH:1:MET:HE2	8:AH:2:LEU:N	2.10	0.66
9:AI:113:LYS:HB2	9:AI:116:LYS:HG3	1.77	0.66
10:AJ:40:LEU:HB2	10:AJ:41:PRO:HD2	1.78	0.66
16:AP:38:TYR:O	16:AP:49:LEU:HA	1.96	0.66
20:AT:50:GLU:HG3	20:AT:51:GLU:N	2.11	0.66
35:BA:2115:G:H4'	35:BA:2166:G:N2	2.10	0.66
35:BA:2199:A:H5'	35:BA:2200:C:OP2	1.96	0.66
35:BA:1758:G:N7	35:BA:2695:C:H4'	2.12	0.66
35:BA:661:C:H4'	46:BP:18:ARG:HG2	1.77	0.66
35:BA:863:A:C2'	35:BA:864:G:H5'	2.26	0.66
36:BB:15:A:H3'	36:BB:16:G:C5'	2.26	0.66
38:BD:4:LYS:NZ	38:BD:20:ASP:HA	2.11	0.66
35:BA:782:A:H2	38:BD:226:MET:HE2	1.60	0.66
41:BG:16:ARG:O	41:BG:20:ILE:HG12	1.96	0.66
44:BN:10:GLU:OE2	44:BN:11:PRO:HD2	1.96	0.66
46:BP:122:PRO:HG3	46:BP:141:ALA:HB3	1.78	0.66
54:BX:83:VAL:O	54:BX:85:PRO:HD3	1.96	0.66
56:BZ:29:TYR:O	56:BZ:30:ASN:HB3	1.96	0.66
1:CA:448:A:O2'	1:CA:449:C:H5'	1.95	0.66
1:CA:601:C:H2'	1:CA:602:A:H8	1.60	0.66
1:CA:990:C:H2'	1:CA:991:U:C6	2.31	0.66
2:CB:70:PHE:O	2:CB:92:TYR:HB2	1.96	0.66
4:CD:101:LEU:HD23	4:CD:121:VAL:HG13	1.78	0.66
6:CF:26:ILE:O	6:CF:30:LEU:HG	1.96	0.66
13:CM:116:THR:HG22	13:CM:117:VAL:N	2.11	0.66
26:DO:2:ALA:O	35:DA:2494:G:H5'	1.96	0.66
29:D3:4:LEU:HD21	29:D3:56:VAL:CG1	2.26	0.66
32:D6:13:CYS:HB3	32:D6:49:HIS:HB3	1.78	0.66
35:DA:2079:U:H3	35:DA:2241:A:H61	1.43	0.66
35:DA:2261:C:H1'	35:DA:2388:A:N3	2.11	0.66
35:DA:449:A:O2'	35:DA:450:G:H5'	1.95	0.66
45:DO:114:ILE:HD12	45:DO:114:ILE:N	2.02	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:104:GLY:O	49:DS:106:ARG:N	2.26	0.66
49:DS:87:PHE:CG	49:DS:88:ASP:N	2.63	0.66
50:DT:22:PHE:H	50:DT:22:PHE:HD2	1.42	0.66
51:DU:31:SER:OG	51:DU:34:LYS:N	2.27	0.66
1:AA:248:C:O2'	1:AA:249:U:H5'	1.96	0.65
2:AB:212:GLN:NE2	2:AB:216:SER:HB2	2.10	0.65
2:AB:48:MET:HA	2:AB:51:LEU:HD12	1.79	0.65
3:AC:111:LEU:HD21	3:AC:145:GLY:O	1.96	0.65
7:AG:49:ILE:HG22	7:AG:49:ILE:O	1.95	0.65
8:AH:39:LEU:O	8:AH:44:PHE:HB2	1.97	0.65
16:AP:56:ALA:O	16:AP:60:LEU:HG	1.95	0.65
34:B8:35:GLN:HA	35:BA:2420:C:P	2.35	0.65
35:BA:1935:G:H3'	35:BA:1962:C:H42	1.62	0.65
35:BA:2119:A:H3'	35:BA:2120:G:C5'	2.26	0.65
35:BA:2208:A:H1'	35:BA:2219:G:C4	2.31	0.65
35:BA:2710:C:OP1	48:BR:15:SER:HB2	1.96	0.65
35:BA:2745:C:H2'	35:BA:2746:U:H6	1.61	0.65
35:BA:380:U:H2'	35:BA:381:G:H8	1.60	0.65
35:BA:703:U:C2'	35:BA:704:G:H5'	2.25	0.65
41:BG:9:ARG:O	41:BG:13:GLU:HG2	1.96	0.65
41:BG:173:LEU:H	41:BG:173:LEU:HD22	1.60	0.65
42:BH:44:VAL:HG12	42:BH:45:VAL:H	1.59	0.65
45:BO:104:ARG:CZ	45:BO:104:ARG:HB3	2.26	0.65
45:BO:37:ASP:H	45:BO:62:VAL:H	1.41	0.65
50:BT:80:SER:CB	50:BT:81:PRO:CD	2.66	0.65
1:CA:248:C:O2'	1:CA:249:U:H5'	1.96	0.65
1:CA:523:A:N1	12:CL:92:ASP:HB2	2.11	0.65
1:CA:946:A:H2'	1:CA:947:G:C8	2.31	0.65
2:CB:172:ILE:CD1	2:CB:172:ILE:H	2.08	0.65
3:CC:173:VAL:O	3:CC:175:LEU:N	2.29	0.65
4:CD:96:LEU:C	4:CD:98:GLU:H	1.99	0.65
8:CH:2:LEU:O	8:CH:3:THR:HG23	1.96	0.65
27:D1:54:ALA:HB2	27:D1:57:GLU:OE1	1.96	0.65
28:D2:20:GLU:O	28:D2:22:GLU:N	2.29	0.65
35:DA:1188:U:C2'	35:DA:1189:A:H5'	2.26	0.65
35:DA:1323:U:H3	35:DA:1331:A:H61	1.44	0.65
35:DA:1570:A:H2'	35:DA:1571:A:H8	1.58	0.65
35:DA:1642:G:H2'	35:DA:1643:G:H8	1.61	0.65
35:DA:1771:C:H2'	35:DA:1772:G:C8	2.30	0.65
1:CA:1517:G:H1'	35:DA:1919:A:O3'	1.96	0.65
35:DA:2682:U:O4	35:DA:2728:U:H1'	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2830:G:C5'	39:DE:58:ARG:HH22	2.09	0.65
40:DF:202:PHE:HE1	40:DF:206:ILE:HD13	1.61	0.65
35:DA:271(P):C:C5'	43:DI:46:ALA:HB2	2.20	0.65
45:DO:17:ARG:O	45:DO:18:LYS:HG3	1.97	0.65
45:DO:32:TYR:CD1	45:DO:32:TYR:N	2.61	0.65
46:DP:74:GLU:OE2	46:DP:75:ILE:HD12	1.96	0.65
50:DT:32:TYR:O	50:DT:33:LYS:HB2	1.95	0.65
50:DT:48:ILE:C	50:DT:63:VAL:HG12	2.16	0.65
1:AA:1048:G:H4'	14:AN:2:ALA:N	2.11	0.65
1:AA:154:C:H2'	1:AA:155:C:H6	1.60	0.65
1:AA:452:A:O2'	1:AA:453:A:H8	1.73	0.65
2:AB:158:LEU:N	2:AB:158:LEU:HD12	2.11	0.65
3:AC:53:ALA:O	3:AC:54:ARG:HB2	1.95	0.65
7:AG:14:PRO:HG3	7:AG:21:VAL:HG13	1.76	0.65
7:AG:86:GLN:NE2	23:AW:32:G:H21	1.93	0.65
10:AJ:48:THR:HG22	10:AJ:49:VAL:N	2.11	0.65
13:AM:9:ILE:HD12	13:AM:9:ILE:N	2.12	0.65
17:AQ:68:ARG:N	17:AQ:70:ARG:HH12	1.95	0.65
35:BA:1210:A:H4'	35:BA:1211:U:O5'	1.95	0.65
39:BE:24:THR:CG2	39:BE:184:VAL:HG23	2.25	0.65
41:BG:76:SER:CB	41:BG:83:ARG:HB3	2.25	0.65
42:BH:156:ALA:C	42:BH:158:HIS:N	2.47	0.65
46:BP:126:VAL:HA	46:BP:145:PRO:HG2	1.78	0.65
47:BQ:51:ARG:O	47:BQ:54:MET:HB3	1.95	0.65
48:BR:18:LEU:HD11	48:BR:22:ARG:CZ	2.26	0.65
51:BU:47:TYR:HD1	51:BU:50:ARG:HH22	1.41	0.65
51:BU:92:ARG:HD2	52:BV:11:GLN:HG3	1.78	0.65
53:BW:6:ILE:HG13	53:BW:104:THR:HG23	1.77	0.65
53:BW:58:ALA:HB1	53:BW:64:MET:SD	2.37	0.65
54:BX:59:VAL:HG23	54:BX:74:PRO:HD2	1.78	0.65
56:BZ:102:LEU:HD11	56:BZ:124:ILE:CG2	2.25	0.65
56:BZ:58:VAL:HG13	56:BZ:67:LEU:C	2.17	0.65
1:CA:1427:U:H2'	1:CA:1428:A:C8	2.30	0.65
1:CA:1399:C:C2	1:CA:1502:A:N6	2.64	0.65
1:CA:331:G:OP1	1:CA:332:G:H5'	1.96	0.65
1:CA:927:G:H2'	1:CA:928:G:C8	2.31	0.65
2:CB:108:ILE:HD13	2:CB:108:ILE:O	1.95	0.65
2:CB:72:GLY:HA3	2:CB:165:VAL:CG1	2.26	0.65
9:CI:28:VAL:HG13	9:CI:64:THR:CA	2.26	0.65
9:CI:70:LYS:O	9:CI:74:ILE:HG13	1.95	0.65
13:CM:49:THR:HG22	13:CM:51:ALA:H	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CP:56:ALA:O	16:CP:60:LEU:HG	1.95	0.65
18:CR:43:PHE:C	18:CR:44:LEU:HD12	2.16	0.65
25:CY:165:THR:HG22	25:CY:166:ASP:N	2.11	0.65
34:D8:6:THR:HG21	34:D8:63:PRO:HD3	1.78	0.65
35:DA:1889:A:O2'	35:DA:2087:G:H5'	1.96	0.65
35:DA:2223:G:C2'	35:DA:2224:G:H5'	2.26	0.65
35:DA:2830:G:H5'	39:DE:58:ARG:NH1	2.12	0.65
35:DA:997:G:O2'	35:DA:998:C:H5'	1.96	0.65
37:DC:47:LEU:HA	37:DC:207:THR:HA	1.78	0.65
38:DD:125:ILE:N	38:DD:125:ILE:HD12	2.12	0.65
40:DF:158:THR:HG21	40:DF:163:VAL:CB	2.26	0.65
43:DI:110:ASP:O	43:DI:112:LYS:N	2.28	0.65
44:DN:30:ILE:O	44:DN:34:LEU:HB2	1.95	0.65
49:DS:26:LEU:HD22	49:DS:87:PHE:CE1	2.31	0.65
45:DO:119:PRO:HB2	50:DT:68:TYR:CD1	2.31	0.65
54:DX:80:ILE:O	54:DX:81:VAL:HB	1.96	0.65
1:AA:1106:G:H2'	1:AA:1107:C:C6	2.31	0.65
1:AA:1513:A:H2'	1:AA:1514:C:H6	1.60	0.65
2:AB:108:ILE:HD13	2:AB:108:ILE:O	1.94	0.65
5:AE:78:HIS:CE1	5:AE:143:ARG:H	2.13	0.65
7:AG:42:ILE:HA	7:AG:45:ASP:HB2	1.78	0.65
7:AG:84:ASN:N	7:AG:84:ASN:HD22	1.92	0.65
9:AI:47:LEU:HB3	9:AI:50:LEU:HD12	1.79	0.65
17:AQ:60:ILE:HG12	17:AQ:61:GLU:O	1.97	0.65
19:AS:22:LEU:HD13	19:AS:22:LEU:O	1.97	0.65
27:B1:77:ALA:C	27:B1:78:LYS:HD2	2.17	0.65
28:B2:20:GLU:O	28:B2:23:LYS:N	2.29	0.65
28:B2:55:ARG:NH2	54:BX:3:THR:HG23	2.11	0.65
35:BA:1697:G:H3'	35:BA:1698:A:C5'	2.26	0.65
35:BA:1999:C:H2'	35:BA:2000:G:H8	1.59	0.65
35:BA:2387:U:H5'	35:BA:2388:A:OP2	1.96	0.65
35:BA:2884:U:H2'	35:BA:2885:C:H5'	1.78	0.65
38:BD:76:PRO:HG2	38:BD:98:VAL:CG2	2.26	0.65
38:BD:80:ALA:HB3	38:BD:94:LEU:HD13	1.77	0.65
41:BG:57:ALA:HB1	41:BG:68:PRO:HB3	1.77	0.65
42:BH:145:ALA:HB1	42:BH:164:TYR:CE1	2.31	0.65
43:BI:37:VAL:HG13	43:BI:38:LEU:CD1	2.26	0.65
43:BI:38:LEU:H	43:BI:38:LEU:CD1	2.02	0.65
44:BN:16:ILE:O	44:BN:54:VAL:HA	1.96	0.65
45:BO:102:VAL:HB	45:BO:106:LEU:CD1	2.25	0.65
45:BO:32:TYR:N	45:BO:32:TYR:CD1	2.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:111:ARG:HA	46:BP:128:HIS:ND1	2.11	0.65
49:BS:67:ARG:O	49:BS:71:ARG:N	2.30	0.65
1:CA:376:G:OP2	16:CP:67:THR:HG21	1.96	0.65
1:CA:696:A:H2'	1:CA:697:U:C6	2.31	0.65
2:CB:157:ARG:HG2	2:CB:158:LEU:H	1.59	0.65
3:CC:111:LEU:HD21	3:CC:145:GLY:O	1.96	0.65
7:CG:36:LYS:HB2	7:CG:36:LYS:HZ2	1.61	0.65
10:CJ:6:ILE:HA	10:CJ:97:GLU:O	1.96	0.65
11:CK:44:SER:N	11:CK:47:VAL:CG2	2.59	0.65
18:CR:50:ILE:CD1	18:CR:70:ILE:HG21	2.26	0.65
27:D1:46:LEU:HD13	27:D1:48:LYS:HE3	1.79	0.65
27:D1:62:VAL:HG21	27:D1:67:ILE:CA	2.23	0.65
31:D5:20:ARG:HH12	53:DW:15:ARG:CZ	2.09	0.65
34:D8:35:GLN:HA	35:DA:2420:C:P	2.35	0.65
35:DA:265:A:H1'	35:DA:266:G:O4'	1.97	0.65
35:DA:506:G:H4'	35:DA:509:C:O2	1.96	0.65
35:DA:977:G:O2'	35:DA:978:G:H5'	1.97	0.65
41:DG:98:ARG:O	41:DG:101:ILE:HG22	1.96	0.65
42:DH:41:MET:HG3	42:DH:53:GLU:O	1.94	0.65
43:DI:129:THR:OG1	43:DI:135:GLU:HB3	1.96	0.65
46:DP:33:ARG:O	46:DP:34:GLY:C	2.34	0.65
49:DS:67:ARG:O	49:DS:71:ARG:N	2.29	0.65
54:DX:34:ALA:O	54:DX:36:LYS:HG3	1.95	0.65
56:DZ:118:GLN:O	56:DZ:172:ALA:HA	1.97	0.65
56:DZ:18:LEU:HG	56:DZ:23:LYS:HB2	1.79	0.65
56:DZ:56:VAL:HG22	56:DZ:70:LEU:HD11	1.77	0.65
1:AA:1242:C:H2'	1:AA:1243:C:H6	1.61	0.65
6:AF:62:TRP:CZ2	6:AF:64:GLN:HB2	2.32	0.65
6:AF:73:ASN:O	6:AF:76:ALA:HB3	1.96	0.65
11:AK:44:SER:OG	11:AK:47:VAL:HG23	1.95	0.65
16:AP:28:ARG:HG2	16:AP:28:ARG:HH11	1.60	0.65
19:AS:48:THR:HG22	19:AS:61:TYR:CA	2.26	0.65
35:BA:1414:G:H2'	35:BA:1415:U:H6	1.59	0.65
35:BA:1515:G:C2'	35:BA:1516:C:H5'	2.27	0.65
35:BA:696:G:C2'	35:BA:697:C:H5'	2.26	0.65
35:BA:892:G:H1	35:BA:894:C:N4	1.94	0.65
38:BD:35:LYS:HE3	38:BD:65:ILE:N	2.09	0.65
39:BE:39:PRO:HA	39:BE:43:GLY:CA	2.25	0.65
40:BF:46:ARG:HG3	40:BF:48:THR:HG23	1.77	0.65
41:BG:57:ALA:HB1	41:BG:68:PRO:CB	2.26	0.65
46:BP:41:ARG:N	46:BP:41:ARG:HD2	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:35:ILE:HG22	53:BW:36:LEU:HD23	1.77	0.65
53:BW:64:MET:O	53:BW:65:LEU:HB3	1.95	0.65
4:CD:104:VAL:HG21	4:CD:140:VAL:HG21	1.77	0.65
4:CD:201:GLN:HA	4:CD:204:ILE:HD12	1.78	0.65
5:CE:109:ILE:HG22	5:CE:110:LEU:N	2.12	0.65
8:CH:18:ARG:N	8:CH:78:GLN:NE2	2.45	0.65
10:CJ:78:ASN:C	10:CJ:80:LYS:H	2.00	0.65
23:CW:42:C:H2'	23:CW:43:G:C8	2.31	0.65
24:CX:14:U:H2'	24:CX:15:G:C8	2.30	0.65
35:DA:1635:G:H5'	35:DA:1635:G:C8	2.31	0.65
35:DA:1713:U:O2'	35:DA:1714:G:H5'	1.97	0.65
25:CY:130:ARG:HH21	35:DA:1942:C:H2'	1.61	0.65
35:DA:2579:C:O2'	35:DA:2580:U:H5'	1.96	0.65
35:DA:2863:C:H2'	35:DA:2864:G:H5''	1.79	0.65
38:DD:166:GLN:HE21	38:DD:166:GLN:N	1.93	0.65
40:DF:37:VAL:O	40:DF:40:GLN:HB2	1.96	0.65
41:DG:37:VAL:CG2	41:DG:99:MET:HG3	2.25	0.65
43:DI:12:LEU:HD12	43:DI:19:VAL:HG11	1.78	0.65
44:DN:46:VAL:HG11	44:DN:48:MET:HG3	1.79	0.65
45:DO:31:LYS:C	45:DO:32:TYR:CD1	2.69	0.65
47:DQ:50:ALA:HA	47:DQ:124:LYS:HG3	1.79	0.65
50:DT:100:TYR:HD2	50:DT:103:ARG:CZ	2.10	0.65
56:DZ:166:SER:HB2	56:DZ:167:PRO:C	2.17	0.65
1:AA:1423:G:H5'	45:BO:49:ARG:HH21	1.61	0.65
1:AA:585:G:H4'	12:AL:8:ASN:ND2	2.06	0.65
1:AA:741:G:H2'	1:AA:742:G:O4'	1.97	0.65
4:AD:126:ILE:HG22	4:AD:127:THR:N	2.11	0.65
5:AE:76:ILE:HD11	5:AE:142:LEU:HD11	1.77	0.65
18:AR:58:LEU:CD1	18:AR:58:LEU:H	2.06	0.65
31:B5:40:LYS:HE2	31:B5:46:CYS:N	2.11	0.65
35:BA:1221:C:H2'	35:BA:1221(A):C:H6	1.61	0.65
35:BA:1280:G:H3'	35:BA:1281:G:H5''	1.78	0.65
35:BA:1841:U:H2'	35:BA:1842:G:C8	2.32	0.65
35:BA:2262:U:H4'	35:BA:2328:A:C2	2.31	0.65
35:BA:49:A:OP1	35:BA:51:G:H5'	1.97	0.65
44:BN:93:THR:HG23	44:BN:93:THR:O	1.97	0.65
35:BA:2413:G:H21	46:BP:70:GLN:HE21	1.44	0.65
46:BP:7:ARG:O	46:BP:10:PRO:HD3	1.95	0.65
53:BW:95:ILE:O	53:BW:95:ILE:HG13	1.96	0.65
55:BY:37:VAL:O	55:BY:38:ILE:HB	1.95	0.65
1:CA:237:C:H4'	17:CQ:25:ARG:NH1	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:32:A:H2'	1:CA:33:A:H8	1.59	0.65
5:CE:96:PRO:HA	5:CE:117:ASP:OD2	1.97	0.65
8:CH:36:LEU:O	8:CH:39:LEU:HB3	1.95	0.65
20:CT:50:GLU:HG3	20:CT:51:GLU:N	2.12	0.65
23:CW:24:C:H2'	23:CW:25:U:C6	2.31	0.65
35:DA:1171:G:H3'	35:DA:1173:G:O4'	1.96	0.65
35:DA:1335:U:H2'	35:DA:1336:A:C8	2.31	0.65
35:DA:1884:A:H2'	35:DA:1885:A:C5'	2.21	0.65
35:DA:200:U:H2'	35:DA:201:C:H5'	1.79	0.65
35:DA:582:G:H2'	35:DA:583:G:H8	1.59	0.65
35:DA:971:C:H2'	35:DA:972:G:O4'	1.96	0.65
41:DG:106:LEU:C	41:DG:108:ASN:H	1.99	0.65
41:DG:144:ILE:HD12	41:DG:145:THR:H	1.60	0.65
44:DN:10:GLU:OE2	44:DN:11:PRO:HD2	1.96	0.65
45:DO:102:VAL:HB	45:DO:106:LEU:CD1	2.26	0.65
55:DY:16:ALA:C	55:DY:21:LYS:HD2	2.17	0.65
56:DZ:166:SER:HB3	56:DZ:168:GLU:H	1.59	0.65
1:AA:1310:G:O2'	1:AA:1311:G:H5'	1.97	0.65
1:AA:340:U:H2'	1:AA:341:C:H6	1.61	0.65
1:AA:403:C:H2'	1:AA:404:U:C6	2.32	0.65
1:AA:471:G:H2'	1:AA:472:A:C8	2.32	0.65
1:AA:927:G:H2'	1:AA:928:G:C8	2.32	0.65
2:AB:162:ILE:C	2:AB:162:ILE:HD12	2.16	0.65
2:AB:33:TYR:HD1	2:AB:43:ASP:HA	1.59	0.65
4:AD:120:LEU:HB3	4:AD:125:HIS:HB2	1.76	0.65
5:AE:126:ARG:C	5:AE:131:ILE:HD11	2.17	0.65
7:AG:101:LEU:O	7:AG:105:VAL:HG23	1.96	0.65
7:AG:11:GLN:HE21	7:AG:12:LEU:H	1.45	0.65
10:AJ:78:ASN:C	10:AJ:80:LYS:H	1.99	0.65
13:AM:116:THR:HG22	13:AM:117:VAL:N	2.11	0.65
18:AR:35:ARG:C	18:AR:37:VAL:H	1.98	0.65
19:AS:53:ASN:ND2	19:AS:55:LYS:H	1.93	0.65
20:AT:89:ARG:HB2	20:AT:104:LEU:HD11	1.78	0.65
25:AY:86:SER:OG	25:AY:88:LEU:HD13	1.96	0.65
32:B6:16:CYS:SG	32:B6:47:THR:HG21	2.35	0.65
35:BA:104:U:H2'	35:BA:105:C:O4'	1.96	0.65
35:BA:1213:A:H2'	35:BA:1214:A:H8	1.61	0.65
35:BA:2830:G:H5'	39:BE:58:ARG:HH22	1.62	0.65
42:BH:73:ALA:O	42:BH:76:VAL:HB	1.97	0.65
44:BN:66:LYS:O	44:BN:67:LEU:HD23	1.95	0.65
45:BO:23:ARG:O	45:BO:39:ILE:HG13	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:75:TYR:N	53:BW:75:TYR:CD1	2.63	0.65
1:CA:1404:C:H2'	1:CA:1405:G:C8	2.32	0.65
1:CA:54:C:H41	1:CA:352:C:H2'	1.61	0.65
2:CB:132:LYS:O	2:CB:135:GLN:HB2	1.97	0.65
3:CC:11:ARG:HH11	3:CC:11:ARG:HG2	1.62	0.65
7:CG:14:PRO:HG3	7:CG:21:VAL:HG13	1.78	0.65
9:CI:113:LYS:HB2	9:CI:116:LYS:HG3	1.78	0.65
18:CR:36:ASN:ND2	18:CR:39:VAL:HB	2.11	0.65
25:CY:21:LEU:HD11	25:CY:121:TYR:O	1.96	0.65
33:D7:25:PRO:HA	33:D7:28:ARG:NH2	2.12	0.65
35:DA:1213:A:H2'	35:DA:1214:A:H8	1.62	0.65
35:DA:1824:G:O2'	35:DA:1825:A:H5'	1.95	0.65
35:DA:2036:C:H6	35:DA:2036:C:C5'	2.05	0.65
35:DA:2119:A:H3'	35:DA:2120:G:C5'	2.27	0.65
35:DA:2467:C:H2'	35:DA:2468:G:H5'	1.79	0.65
35:DA:2677:G:H2'	35:DA:2678:C:C6	2.30	0.65
35:DA:611:C:O2'	35:DA:612:C:H5'	1.96	0.65
35:DA:933:A:H2'	35:DA:934:G:O4'	1.96	0.65
38:DD:35:LYS:HE3	38:DD:65:ILE:N	2.10	0.65
39:DE:52:LEU:CB	39:DE:76:ARG:HB2	2.26	0.65
40:DF:28:ILE:HG22	40:DF:112:MET:HB3	1.79	0.65
35:DA:323:G:H2'	40:DF:169:ASN:ND2	2.10	0.65
48:DR:18:LEU:HD11	48:DR:22:ARG:CZ	2.27	0.65
1:AA:1166:G:N2	1:AA:1169:A:H3'	2.12	0.65
1:AA:320:C:H2'	1:AA:321:A:C8	2.31	0.65
1:AA:783:C:O2'	1:AA:784:C:H5'	1.96	0.65
2:AB:157:ARG:HG2	2:AB:158:LEU:H	1.61	0.65
9:AI:48:GLU:N	9:AI:49:PRO:HD2	2.12	0.65
15:AO:17:ARG:HH11	15:AO:17:ARG:HG3	1.61	0.65
15:AO:56:LEU:O	15:AO:60:VAL:HG23	1.97	0.65
31:B5:46:CYS:SG	31:B5:47:PRO:HD2	2.37	0.65
33:B7:30:VAL:HG13	33:B7:33:ARG:HH22	1.60	0.65
34:B8:42:ARG:O	34:B8:44:LYS:N	2.29	0.65
35:BA:2415:G:H4'	46:BP:66:GLY:HA3	1.78	0.65
35:BA:2861:G:O2'	35:BA:2862:G:H5'	1.97	0.65
35:BA:2866:U:C5	35:BA:2868:A:H1'	2.31	0.65
35:BA:285:C:C2'	35:BA:286:C:H5"	2.27	0.65
35:BA:970:C:H2'	35:BA:971:C:H6	1.61	0.65
35:BA:999:U:H2'	35:BA:999:U:O2	1.97	0.65
38:BD:94:LEU:O	38:BD:94:LEU:HD13	1.96	0.65
1:CA:1116:C:C3'	1:CA:1117:G:H5"	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:376:G:H5''	16:CP:5:ARG:HD2	1.79	0.65
1:CA:748:C:H1'	1:CA:749:C:OP2	1.95	0.65
11:CK:103:LEU:N	11:CK:103:LEU:HD22	2.06	0.65
12:CL:27:LEU:O	12:CL:29:GLY:N	2.30	0.65
17:CQ:9:VAL:HG12	17:CQ:10:VAL:H	1.62	0.65
18:CR:35:ARG:C	18:CR:37:VAL:H	1.98	0.65
25:CY:133:ARG:HG3	25:CY:161:ILE:HG22	1.77	0.65
26:D0:77:ARG:NH2	35:DA:857:C:H5'	2.11	0.65
35:DA:1018:C:O2'	35:DA:1019:U:H5'	1.96	0.65
35:DA:2074:U:H2'	35:DA:2075:U:C6	2.32	0.65
35:DA:783:A:H2'	35:DA:784:A:H4'	1.77	0.65
42:DH:89:ILE:CD1	42:DH:89:ILE:N	2.59	0.65
45:DO:104:ARG:CZ	45:DO:104:ARG:HB3	2.27	0.65
45:DO:31:LYS:HD2	45:DO:32:TYR:HE1	1.61	0.65
48:DR:82:GLU:C	48:DR:85:PRO:HD2	2.16	0.65
52:DV:19:LYS:HE2	52:DV:19:LYS:HA	1.79	0.65
35:DA:975(A):G:OP1	52:DV:79:VAL:HG13	1.96	0.65
53:DW:6:ILE:HG13	53:DW:104:THR:HG23	1.77	0.65
56:DZ:109:ALA:HB3	56:DZ:145:GLU:OE1	1.97	0.65
1:AA:112:G:H4'	1:AA:389:A:H5''	1.78	0.65
1:AA:192:U:C4'	20:AT:103:GLY:HA2	2.26	0.65
1:AA:254:G:OP1	17:AQ:67:LYS:O	2.15	0.65
3:AC:164:ARG:HB2	3:AC:164:ARG:NH1	2.07	0.65
4:AD:13:ARG:HG2	4:AD:14:ARG:N	2.12	0.65
8:AH:82:HIS:CD2	8:AH:138:TRP:HE1	2.09	0.65
9:AI:7:THR:HB	9:AI:83:ARG:NH1	2.12	0.65
14:AN:4:LYS:O	14:AN:7:ILE:HG12	1.97	0.65
15:AO:30:ALA:O	15:AO:33:THR:HB	1.97	0.65
19:AS:62:ILE:HD12	19:AS:66:MET:HG3	1.79	0.65
20:AT:63:ILE:HD12	20:AT:81:LYS:HG2	1.79	0.65
25:AY:174:GLN:O	25:AY:177:GLU:HB3	1.97	0.65
25:AY:42:LYS:HB3	25:AY:49:HIS:HB3	1.79	0.65
31:B5:49:CYS:O	31:B5:56:LYS:HB2	1.97	0.65
35:BA:1366:A:H2'	35:BA:1367:A:H8	1.62	0.65
35:BA:1406:U:H3'	35:BA:1407:C:H6	1.62	0.65
35:BA:1587:A:H2'	35:BA:1588:C:O4'	1.97	0.65
35:BA:1681:G:OP2	35:BA:1681:G:H8	1.80	0.65
35:BA:2111:C:O2'	35:BA:2118:U:H4'	1.97	0.65
35:BA:225:A:O2'	35:BA:226:G:H5'	1.96	0.65
35:BA:2863:C:H2'	35:BA:2864:G:H5''	1.79	0.65
44:BN:3:THR:HG22	44:BN:5:VAL:CG2	2.25	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:22:PHE:HD2	50:BT:22:PHE:H	1.44	0.65
54:BX:77:LYS:HE3	54:BX:78:LYS:N	2.12	0.65
55:BY:31:LEU:HD23	55:BY:36:ALA:HB3	1.77	0.65
1:CA:924:C:H2'	1:CA:925:G:C8	2.32	0.65
3:CC:6:HIS:NE2	3:CC:8:ILE:HB	2.11	0.65
4:CD:200:GLU:O	4:CD:204:ILE:HG13	1.96	0.65
4:CD:96:LEU:N	4:CD:96:LEU:HD22	2.05	0.65
8:CH:82:HIS:CD2	8:CH:138:TRP:HE1	2.09	0.65
8:CH:22:GLU:O	8:CH:63:LEU:HD23	1.96	0.65
11:CK:86:GLY:H	11:CK:112:THR:HG23	1.61	0.65
12:CL:34:ARG:HB3	12:CL:61:THR:HG21	1.77	0.65
19:CS:48:THR:HG22	19:CS:61:TYR:CA	2.26	0.65
20:CT:89:ARG:HB2	20:CT:104:LEU:HD11	1.79	0.65
23:CW:35:C:H2'	23:CW:36:A:O4'	1.96	0.65
28:D2:49:LYS:HE2	35:DA:76:C:OP1	1.97	0.65
34:D8:43:GLN:C	34:D8:44:LYS:HD2	2.16	0.65
35:DA:1854:A:H3'	35:DA:1855:G:H8	1.61	0.65
35:DA:2410:G:C2	35:DA:2411:A:H1'	2.31	0.65
35:DA:2506:U:H4'	35:DA:2507:C:OP1	1.95	0.65
35:DA:393:C:H2'	35:DA:394:A:H8	1.61	0.65
26:D0:26:TYR:CE2	35:DA:857:C:H1'	2.32	0.65
38:DD:268:ARG:HH11	38:DD:268:ARG:HB2	1.62	0.65
40:DF:2:LYS:O	40:DF:25:PRO:HD2	1.97	0.65
42:DH:116:GLU:HG2	42:DH:117:PRO:HD2	1.77	0.65
43:DI:5:LEU:N	43:DI:5:LEU:HD23	2.12	0.65
44:DN:56:ASN:CA	44:DN:124:ALA:HA	2.27	0.65
44:DN:93:THR:O	44:DN:93:THR:HG23	1.97	0.65
28:D2:26:ARG:HD3	54:DX:5:TYR:HD1	1.61	0.65
54:DX:59:VAL:C	54:DX:73:ARG:HA	2.18	0.65
54:DX:77:LYS:HE2	54:DX:78:LYS:HG3	1.77	0.65
55:DY:28:LYS:HD2	55:DY:37:VAL:CG1	2.26	0.65
1:AA:104:G:O2'	1:AA:105:G:H5'	1.97	0.65
1:AA:884:U:H4'	1:AA:885:G:H5''	1.79	0.65
2:AB:31:TYR:CD2	2:AB:31:TYR:N	2.60	0.65
4:AD:191:ARG:HH12	4:AD:195:ALA:HA	1.61	0.65
12:AL:34:ARG:HB3	12:AL:61:THR:HG21	1.79	0.65
13:AM:65:LYS:C	13:AM:66:LEU:N	2.50	0.65
15:AO:28:GLN:O	15:AO:32:LEU:HG	1.97	0.65
21:AU:8:THR:O	21:AU:12:LYS:HB2	1.97	0.65
25:AY:148:HIS:O	25:AY:149:LEU:HD23	1.95	0.65
25:AY:43:VAL:HG13	25:AY:82:ALA:HB3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:51:VAL:HG23	27:B1:62:VAL:CG1	2.26	0.65
31:B5:17:ASP:O	31:B5:20:ARG:HB2	1.96	0.65
35:BA:1742:G:N7	35:BA:1743:C:N3	2.44	0.65
35:BA:1765:C:H2'	35:BA:1766:U:C6	2.31	0.65
35:BA:2292:C:O2'	35:BA:2293:C:H5'	1.95	0.65
35:BA:967:C:O2'	35:BA:968:G:H5'	1.97	0.65
38:BD:92:ILE:HA	38:BD:107:ALA:H	1.60	0.65
39:BE:4:ILE:CG1	39:BE:28:ALA:HB1	2.27	0.65
41:BG:15:VAL:O	41:BG:19:LEU:HG	1.97	0.65
46:BP:17:LYS:O	46:BP:19:VAL:N	2.29	0.65
47:BQ:34:LEU:HD12	47:BQ:35:VAL:N	2.11	0.65
56:BZ:10:ARG:HG2	56:BZ:11:GLU:H	1.60	0.65
56:BZ:29:TYR:CE2	56:BZ:87:ASP:HB2	2.31	0.65
56:BZ:71:VAL:HG22	56:BZ:88:PHE:CE2	2.32	0.65
1:CA:1030(A):G:H1'	1:CA:1031:G:H1	1.62	0.65
1:CA:1456:G:C2'	1:CA:1457:G:H5'	2.26	0.65
1:CA:154:C:H2'	1:CA:155:C:H6	1.60	0.65
1:CA:253:U:H2'	1:CA:254:G:C8	2.32	0.65
3:CC:53:ALA:O	3:CC:54:ARG:HB2	1.94	0.65
4:CD:192:GLU:H	4:CD:192:GLU:CD	2.01	0.65
7:CG:40:ALA:HA	7:CG:43:PHE:HB3	1.79	0.65
25:CY:169:ILE:O	25:CY:170:ALA:C	2.35	0.65
29:D3:15:TYR:HB3	29:D3:19:GLN:NE2	2.12	0.65
35:DA:1937:A:C2'	35:DA:1938:A:H5'	2.26	0.65
35:DA:554:U:O2'	35:DA:555:U:H5'	1.96	0.65
35:DA:816:C:O2'	35:DA:817:C:H5'	1.97	0.65
38:DD:68:LYS:O	38:DD:68:LYS:HG3	1.96	0.65
35:DA:2051:A:H4'	39:DE:141:ILE:HD11	1.79	0.65
39:DE:2:LYS:HE2	39:DE:95:ILE:HG22	1.79	0.65
47:DQ:64:ILE:HG23	47:DQ:106:VAL:HG13	1.77	0.65
49:DS:24:LEU:HB2	49:DS:85:VAL:HB	1.79	0.65
49:DS:25:ARG:HH21	49:DS:89:ARG:NH1	1.93	0.65
51:DU:3:ARG:HH11	51:DU:3:ARG:CG	2.08	0.65
35:DA:993:G:H5'	52:DV:75:PHE:CZ	2.32	0.65
56:DZ:132:ASN:O	56:DZ:134:PRO:HD3	1.96	0.65
1:AA:1118:C:OP1	9:AI:9:ARG:HD3	1.97	0.65
1:AA:1203:C:H2'	1:AA:1204:A:H8	1.62	0.65
2:AB:137:ARG:HG2	2:AB:137:ARG:HH11	1.61	0.65
4:AD:192:GLU:CD	4:AD:192:GLU:H	2.01	0.65
4:AD:65:ARG:HH11	4:AD:72:GLU:N	1.95	0.65
9:AI:77:ILE:O	9:AI:81:ILE:HG12	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:85:ILE:HD12	12:AL:99:HIS:O	1.96	0.65
26:B0:2:ALA:O	35:BA:2494:G:H5'	1.97	0.65
26:B0:72:ARG:HD3	26:B0:75:LEU:HD13	1.77	0.65
35:BA:1568:G:H5''	38:BD:61:LEU:HB2	1.79	0.65
35:BA:1819:A:H1'	35:BA:1821:A:C6	2.32	0.65
35:BA:2487:G:H2'	35:BA:2488:A:C8	2.31	0.65
35:BA:2051:A:H5'	35:BA:2578:G:O4'	1.96	0.65
35:BA:2795:G:H2'	35:BA:2795:G:N3	2.10	0.65
35:BA:882:G:H22	35:BA:895:U:H3	1.43	0.65
35:BA:997:G:O2'	35:BA:998:C:H5'	1.97	0.65
41:BG:11:TYR:O	41:BG:15:VAL:HB	1.97	0.65
41:BG:36:LYS:CE	41:BG:160:VAL:HG21	2.26	0.65
42:BH:145:ALA:HB1	42:BH:164:TYR:HE1	1.62	0.65
43:BI:51:ILE:HG22	43:BI:52:ARG:N	2.12	0.65
54:BX:33:LYS:C	54:BX:35:THR:N	2.46	0.65
3:CC:16:ARG:HH11	3:CC:16:ARG:HA	1.60	0.65
3:CC:182:ILE:HG23	3:CC:203:PHE:HA	1.78	0.65
8:CH:12:ARG:NH1	8:CH:26:VAL:HA	2.12	0.65
31:D5:44:THR:HG21	48:DR:101:ALA:CB	2.23	0.65
31:D5:49:CYS:HB2	31:D5:59:GLU:OE1	1.96	0.65
35:DA:1681:G:OP2	35:DA:1681:G:H8	1.80	0.65
35:DA:2206:G:C2	35:DA:2207:G:H5'	2.31	0.65
35:DA:2290:G:H8	35:DA:2290:G:H5'	1.61	0.65
35:DA:2884:U:H2'	35:DA:2885:C:H5'	1.78	0.65
35:DA:909:A:H1'	47:DQ:10:ARG:NH2	2.12	0.65
36:DB:20:C:C2'	36:DB:21:G:H5''	2.26	0.65
40:DF:60:SER:OG	40:DF:61:GLY:N	2.28	0.65
49:DS:13:ARG:H	49:DS:13:ARG:CD	1.99	0.65
51:DU:92:ARG:C	51:DU:94:ASN:H	1.99	0.65
56:DZ:53:ILE:HG13	56:DZ:53:ILE:O	1.95	0.65
1:AA:1030:C:H2'	1:AA:1030(A):G:H5'	1.77	0.64
1:AA:586:C:O2'	1:AA:587:G:H5'	1.97	0.64
1:AA:642:A:N3	8:AH:113:SER:OG	2.24	0.64
3:AC:16:ARG:HA	3:AC:16:ARG:HH11	1.62	0.64
11:AK:44:SER:N	11:AK:47:VAL:CG2	2.61	0.64
12:AL:37:CYS:SG	12:AL:81:SER:HB2	2.36	0.64
20:AT:31:SER:O	20:AT:34:LYS:HB2	1.97	0.64
23:AW:39:A:H2'	23:AW:40:C:H5'	1.79	0.64
27:B1:20:ARG:NH1	27:B1:41:ARG:NE	2.45	0.64
35:BA:2591:C:H2'	35:BA:2592:G:H8	1.58	0.64
35:BA:678:C:H2'	35:BA:679:C:C6	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1353:A:H4'	38:BD:38:LYS:NZ	2.12	0.64
42:BH:44:VAL:HG12	42:BH:45:VAL:N	2.12	0.64
42:BH:89:ILE:CD1	42:BH:90:LYS:H	2.10	0.64
45:BO:69:ILE:CD1	45:BO:77:ILE:HG23	2.27	0.64
52:BV:52:VAL:C	52:BV:54:GLY:H	1.99	0.64
1:CA:580:U:H2'	1:CA:581:G:O4'	1.97	0.64
1:CA:734:G:O2'	1:CA:735:C:H5'	1.97	0.64
1:CA:884:U:H4'	1:CA:885:G:H5''	1.78	0.64
4:CD:145:GLU:HG2	4:CD:184:LYS:CG	2.25	0.64
6:CF:5:GLU:HG3	6:CF:93:SER:OG	1.96	0.64
8:CH:10:LEU:HD23	8:CH:10:LEU:H	1.62	0.64
18:CR:19:LYS:O	18:CR:20:ALA:HB2	1.97	0.64
18:CR:44:LEU:O	18:CR:45:SER:HB3	1.96	0.64
18:CR:53:ARG:NH2	18:CR:60:ALA:N	2.44	0.64
25:CY:123:GLU:O	25:CY:126:ARG:N	2.30	0.64
25:CY:32:ARG:HB3	25:CY:103:ILE:HD13	1.80	0.64
25:CY:70:SER:CB	25:CY:76:LEU:HB2	2.27	0.64
27:D1:42:GLN:HG2	27:D1:43:TYR:H	1.61	0.64
29:D3:4:LEU:HD23	29:D3:5:LYS:N	2.13	0.64
35:DA:2014:A:H4'	53:DW:94:ASP:OD1	1.97	0.64
35:DA:2186:G:C3'	35:DA:2187:G:H5''	2.27	0.64
35:DA:2075:U:H2'	35:DA:2238:G:N2	2.12	0.64
38:DD:109:ASP:HB3	38:DD:195:ALA:HB3	1.79	0.64
35:DA:2631:G:N2	39:DE:61:ARG:NH1	2.45	0.64
40:DF:134:GLY:H	40:DF:162:LEU:HD11	1.61	0.64
35:DA:322:A:P	40:DF:169:ASN:HB2	2.37	0.64
40:DF:96:ASP:OD1	40:DF:98:SER:HB3	1.97	0.64
41:DG:114:ILE:HG21	41:DG:117:PHE:CB	2.17	0.64
41:DG:132:ASN:ND2	41:DG:133:LEU:N	2.45	0.64
41:DG:60:LEU:O	41:DG:64:THR:HG22	1.97	0.64
41:DG:73:ALA:H	41:DG:87:PRO:HD2	1.62	0.64
41:DG:4:ASP:HB3	41:DG:8:LYS:CE	2.27	0.64
41:DG:92:VAL:HG22	41:DG:93:THR:H	1.61	0.64
45:DO:43:VAL:HG21	45:DO:52:VAL:HG12	1.78	0.64
45:DO:47:ILE:HG23	45:DO:48:PRO:HD2	1.78	0.64
47:DQ:82:ARG:HG2	47:DQ:82:ARG:HH11	1.61	0.64
51:DU:59:ARG:O	51:DU:61:TRP:N	2.30	0.64
52:DV:39:LEU:HD21	52:DV:53:GLU:HA	1.79	0.64
1:AA:556:C:O2'	1:AA:557:G:H5'	1.97	0.64
2:AB:165:VAL:CG2	2:AB:166:ASP:H	1.99	0.64
3:AC:110:ASN:O	3:AC:141:VAL:HG13	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:182:ILE:HG23	3:AC:202:ILE:C	2.17	0.64
6:AF:5:GLU:HG3	6:AF:93:SER:OG	1.97	0.64
7:AG:113:GLU:CB	7:AG:119:ARG:HG2	2.28	0.64
9:AI:111:ARG:HG2	9:AI:112:LYS:N	2.12	0.64
25:AY:108:GLU:HA	25:AY:111:ARG:CG	2.28	0.64
26:B0:40:GLN:HE21	26:B0:43:THR:HA	1.62	0.64
34:B8:22:VAL:HG21	34:B8:56:GLU:HB2	1.80	0.64
34:B8:50:LEU:C	34:B8:53:PRO:HD2	2.17	0.64
35:BA:2404:C:H2'	35:BA:2405:G:O4'	1.98	0.64
35:BA:2673:G:H2'	35:BA:2674:G:H8	1.62	0.64
35:BA:690:G:H2'	35:BA:691:C:C6	2.32	0.64
35:BA:93:G:H2'	35:BA:94:C:H6	1.62	0.64
38:BD:117:VAL:HG22	38:BD:118:VAL:H	1.61	0.64
39:BE:133:LYS:C	39:BE:134:ILE:HD13	2.17	0.64
40:BF:114:VAL:HG11	40:BF:202:PHE:HE2	1.61	0.64
44:BN:42:TRP:CE2	44:BN:44:PRO:HD3	2.32	0.64
45:BO:36:GLY:H	45:BO:62:VAL:HB	1.59	0.64
48:BR:18:LEU:HD11	48:BR:22:ARG:NH2	2.12	0.64
48:BR:97:VAL:HG22	48:BR:114:VAL:HG22	1.80	0.64
50:BT:107:ASP:OD1	50:BT:109:GLU:HB2	1.97	0.64
50:BT:27:THR:C	50:BT:88:ILE:HD13	2.17	0.64
50:BT:28:VAL:HG11	50:BT:46:GLU:HA	1.78	0.64
54:BX:65:ARG:NH2	54:BX:66:LEU:H	1.95	0.64
55:BY:31:LEU:CB	55:BY:36:ALA:H	2.10	0.64
56:BZ:117:LEU:HA	56:BZ:173:ALA:O	1.96	0.64
1:CA:1310:G:O2'	1:CA:1311:G:H5'	1.96	0.64
1:CA:1381:U:H2'	1:CA:1382:C:H5'	1.79	0.64
2:CB:185:ILE:CG2	2:CB:199:TYR:HB2	2.26	0.64
9:CI:105:ASP:HB3	9:CI:107:ARG:HG3	1.79	0.64
9:CI:95:LYS:HZ3	9:CI:96:LEU:HB2	1.62	0.64
25:CY:171:LYS:HD2	25:CY:175:LEU:HD13	1.77	0.64
32:D6:16:CYS:SG	32:D6:47:THR:HG21	2.37	0.64
34:D8:50:LEU:C	34:D8:53:PRO:HD2	2.18	0.64
35:DA:1159:U:H2'	35:DA:1160:G:H5'	1.78	0.64
35:DA:1353:A:H4'	38:DD:38:LYS:NZ	2.13	0.64
35:DA:1879:C:C3'	35:DA:1880:C:H5''	2.27	0.64
35:DA:2013:A:H4'	53:DW:96:ILE:HD12	1.77	0.64
35:DA:2111:C:O2'	35:DA:2118:U:H4'	1.97	0.64
35:DA:2222:G:O2'	35:DA:2223:G:H5'	1.97	0.64
35:DA:2680:C:H2'	35:DA:2681:C:O2	1.97	0.64
35:DA:2723:C:O2'	35:DA:2724:C:H5'	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:528:A:C2	35:DA:2042:A:H2'	2.32	0.64
38:DD:142:VAL:HA	38:DD:194:GLY:H	1.62	0.64
38:DD:76:PRO:HG2	38:DD:98:VAL:CG2	2.27	0.64
40:DF:22:ALA:O	40:DF:26:ALA:HB2	1.96	0.64
40:DF:5:ALA:O	40:DF:6:VAL:HG13	1.97	0.64
40:DF:88:VAL:HG22	40:DF:89:VAL:N	2.12	0.64
35:DA:2304:G:H4'	41:DG:133:LEU:HB3	1.78	0.64
35:DA:2009:G:H1'	48:DR:107:ASP:O	1.98	0.64
50:DT:109:GLU:HA	50:DT:112:ARG:HG3	1.78	0.64
53:DW:1:MET:HG3	53:DW:2:GLU:H	1.62	0.64
1:AA:1422:G:H2'	1:AA:1423:G:H8	1.62	0.64
1:AA:312:C:H2'	1:AA:313:A:H8	1.62	0.64
1:AA:59:A:H5''	1:AA:60:A:H5''	1.78	0.64
1:AA:880:C:O2'	1:AA:881:G:H5'	1.96	0.64
12:AL:25:PRO:HD2	12:AL:98:TYR:OH	1.96	0.64
16:AP:6:LEU:CD1	16:AP:6:LEU:N	2.60	0.64
19:AS:36:ARG:HH12	19:AS:75:ALA:HB3	1.63	0.64
27:B1:68:PRO:HG2	27:B1:69:LYS:N	2.09	0.64
33:B7:25:PRO:HG2	33:B7:26:GLY:H	1.61	0.64
35:BA:142(A):C:H2'	35:BA:143:G:O4'	1.98	0.64
35:BA:2106:G:H2'	35:BA:2107:C:O4'	1.97	0.64
35:BA:230:U:H2'	35:BA:230:U:O2	1.96	0.64
35:BA:519:U:H5''	53:BW:25:ARG:NH2	2.13	0.64
38:BD:177:LEU:HD12	38:BD:181:GLU:CG	2.27	0.64
40:BF:141:ALA:O	40:BF:144:LYS:HB3	1.98	0.64
40:BF:202:PHE:HE1	40:BF:206:ILE:HD13	1.63	0.64
42:BH:102:ALA:HB2	42:BH:117:PRO:CD	2.19	0.64
43:BI:28:ASN:O	43:BI:32:PRO:HG2	1.98	0.64
43:BI:92:VAL:O	43:BI:92:VAL:HG22	1.96	0.64
49:BS:54:LEU:HD21	49:BS:59:LYS:O	1.98	0.64
50:BT:42:ILE:O	50:BT:42:ILE:HG13	1.97	0.64
56:BZ:149:SER:OG	56:BZ:173:ALA:HB2	1.98	0.64
1:CA:1166:G:N2	1:CA:1169:A:H3'	2.11	0.64
1:CA:174:C:O2'	1:CA:175:C:H5'	1.97	0.64
1:CA:639:G:O2'	1:CA:640:A:H5'	1.98	0.64
3:CC:16:ARG:HH11	3:CC:16:ARG:CB	2.10	0.64
3:CC:182:ILE:HG23	3:CC:202:ILE:C	2.17	0.64
6:CF:30:LEU:HD23	6:CF:75:LEU:HD21	1.79	0.64
9:CI:113:LYS:HD2	9:CI:113:LYS:N	2.13	0.64
16:CP:6:LEU:HB3	16:CP:17:TYR:CD2	2.33	0.64
17:CQ:29:HIS:HB3	17:CQ:33:GLY:N	2.11	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CT:36:LEU:H	20:CT:36:LEU:CD2	2.11	0.64
26:D0:25:ARG:HB2	26:D0:37:LEU:HD23	1.78	0.64
27:D1:34:THR:HG21	35:DA:387:U:O3'	1.97	0.64
35:DA:1605:C:H5'	35:DA:1610:A:N6	2.13	0.64
35:DA:1795:C:H42	35:DA:1824:G:H1	1.45	0.64
42:DH:31:GLY:O	42:DH:79:VAL:HG11	1.97	0.64
49:DS:38:GLN:C	49:DS:39:ILE:HD12	2.18	0.64
51:DU:47:TYR:HA	51:DU:50:ARG:CZ	2.28	0.64
35:DA:1614:A:N6	53:DW:93:ALA:CB	2.60	0.64
54:DX:82:GLN:CG	54:DX:83:VAL:H	2.10	0.64
1:AA:1163:C:H2'	1:AA:1164:G:H8	1.61	0.64
1:AA:1218:C:H2'	1:AA:1219:U:C6	2.32	0.64
1:AA:233:C:H2'	1:AA:234:C:H6	1.62	0.64
1:AA:834:C:H2'	1:AA:835:U:C6	2.33	0.64
2:AB:105:PHE:O	2:AB:108:ILE:N	2.31	0.64
3:AC:134:ILE:HD11	3:AC:153:VAL:HG21	1.78	0.64
7:AG:135:VAL:O	7:AG:138:LYS:HB3	1.96	0.64
8:AH:97:VAL:HG13	8:AH:98:LYS:HG3	1.79	0.64
25:AY:139:LYS:O	25:AY:143:LEU:HB2	1.98	0.64
35:BA:1298:C:H2'	35:BA:1299:G:C8	2.31	0.64
35:BA:2250:G:C5	47:BQ:82:ARG:HD2	2.32	0.64
35:BA:2631:G:N2	39:BE:61:ARG:NH1	2.44	0.64
35:BA:2854:G:H2'	35:BA:2855:C:H6	1.63	0.64
35:BA:661:C:H2'	35:BA:662:G:C8	2.32	0.64
35:BA:759:G:H2'	35:BA:760:G:C8	2.31	0.64
36:BB:91:C:H2'	36:BB:92:C:C6	2.32	0.64
41:BG:36:LYS:HE2	41:BG:160:VAL:HG21	1.79	0.64
44:BN:17:ASP:O	44:BN:19:GLU:N	2.30	0.64
45:BO:120:GLU:HG3	45:BO:122:LEU:HD11	1.80	0.64
46:BP:124:LYS:HA	46:BP:143:GLY:N	2.12	0.64
48:BR:2:ARG:CZ	48:BR:5:LYS:HE3	2.27	0.64
48:BR:60:LEU:HD23	48:BR:61:HIS:N	2.11	0.64
50:BT:3:ARG:HB3	50:BT:6:LEU:HB3	1.79	0.64
50:BT:52:ILE:HG22	50:BT:61:PHE:CB	2.27	0.64
52:BV:1:MET:HE3	52:BV:45:THR:H	1.63	0.64
56:BZ:128:VAL:HG22	56:BZ:132:ASN:HB2	1.78	0.64
1:CA:1090:U:H2'	1:CA:1091:U:H6	1.62	0.64
1:CA:1242:C:H2'	1:CA:1243:C:H6	1.61	0.64
1:CA:786:G:H1	1:CA:796:C:H42	1.45	0.64
3:CC:138:VAL:HG22	3:CC:151:VAL:HG23	1.78	0.64
4:CD:127:THR:HG22	4:CD:149:ALA:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:144:THR:O	5:CE:147:ASP:OD2	2.15	0.64
1:CA:1342:C:H1'	9:CI:124:GLN:NE2	2.12	0.64
13:CM:91:ARG:HH11	19:CS:81:ARG:NH2	1.88	0.64
32:D6:32:ASN:ND2	32:D6:33:LYS:H	1.96	0.64
35:DA:1220:A:O2'	35:DA:1221:C:H5''	1.98	0.64
35:DA:1440:G:H2'	35:DA:1441:G:C8	2.32	0.64
35:DA:2312:U:C2'	35:DA:2313:C:H5''	2.28	0.64
35:DA:2415:G:H4'	46:DP:66:GLY:HA3	1.79	0.64
35:DA:608:A:OP1	40:DF:100:THR:HG21	1.96	0.64
39:DE:141:ILE:N	39:DE:141:ILE:HD13	2.12	0.64
39:DE:1:MET:HG2	39:DE:83:ASP:O	1.97	0.64
40:DF:205:ARG:O	40:DF:205:ARG:HG2	1.96	0.64
46:DP:101:VAL:CG2	46:DP:107:LYS:HA	2.24	0.64
46:DP:95:VAL:HG23	46:DP:125:VAL:HB	1.80	0.64
47:DQ:43:THR:HG1	47:DQ:45:GLN:HB2	1.63	0.64
52:DV:61:VAL:HG12	52:DV:62:LEU:H	1.62	0.64
1:AA:1045:C:H2'	1:AA:1046:A:O4'	1.97	0.64
1:AA:186:C:H4'	20:AT:82:SER:HB3	1.79	0.64
1:AA:337:C:H2'	1:AA:338:A:C8	2.33	0.64
2:AB:178:ARG:HB2	2:AB:178:ARG:HH11	1.62	0.64
4:AD:17:VAL:HG12	4:AD:18:LYS:N	2.11	0.64
4:AD:18:LYS:NZ	4:AD:31:CYS:SG	2.66	0.64
1:AA:426:G:H4'	4:AD:41:GLY:O	1.98	0.64
13:AM:49:THR:HG22	13:AM:51:ALA:H	1.62	0.64
3:AC:9:GLY:HA3	14:AN:49:HIS:HA	1.79	0.64
17:AQ:95:TYR:C	17:AQ:97:SER:H	2.01	0.64
1:AA:333:G:H4'	20:AT:16:HIS:CE1	2.32	0.64
25:AY:29:ARG:NH2	25:AY:32:ARG:HH22	1.94	0.64
35:BA:1171:G:H3'	35:BA:1173:G:O4'	1.96	0.64
35:BA:1602:U:H3'	35:BA:1603:A:H5''	1.80	0.64
35:BA:2028:U:H2'	35:BA:2029:G:C8	2.32	0.64
35:BA:2469:A:H2	35:BA:2481:G:H21	1.46	0.64
40:BF:205:ARG:O	40:BF:205:ARG:HG2	1.96	0.64
45:BO:17:ARG:O	45:BO:18:LYS:HG3	1.97	0.64
1:CA:1057:G:O2'	1:CA:1058:G:H5'	1.97	0.64
1:CA:1507:A:C2	1:CA:1530:G:H1'	2.33	0.64
1:CA:403:C:H2'	1:CA:404:U:C6	2.32	0.64
4:CD:59:ARG:HA	4:CD:59:ARG:NH1	2.12	0.64
7:CG:11:GLN:NE2	7:CG:12:LEU:H	1.95	0.64
10:CJ:5:ARG:O	10:CJ:98:ILE:HA	1.96	0.64
10:CJ:32:ALA:HB1	10:CJ:75:ILE:CG1	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CT:57:ARG:HH11	20:CT:57:ARG:HB2	1.63	0.64
35:DA:1841:U:H2'	35:DA:1842:G:C8	2.32	0.64
35:DA:1879:C:H2'	35:DA:1880:C:C5'	2.19	0.64
35:DA:816:C:H2'	35:DA:817:C:H6	1.62	0.64
36:DB:115:G:O4'	49:DS:47:THR:HB	1.98	0.64
39:DE:64:LYS:C	39:DE:66:HIS:H	2.00	0.64
43:DI:91:SER:H	43:DI:121:LYS:CE	2.10	0.64
49:DS:54:LEU:HD13	49:DS:58:LEU:O	1.97	0.64
1:AA:105:G:H2'	1:AA:106:C:C6	2.33	0.64
1:AA:1381:U:H2'	1:AA:1382:C:H5'	1.80	0.64
1:AA:1502:A:H2	1:AA:1505:G:H1	1.45	0.64
2:AB:69:LEU:HD23	2:AB:159:PRO:HG2	1.78	0.64
2:AB:73:THR:HG22	2:AB:94:ASN:HA	1.79	0.64
3:AC:23:TYR:HA	10:AJ:11:PHE:CE1	2.33	0.64
4:AD:187:ARG:HH11	4:AD:187:ARG:HG2	1.63	0.64
10:AJ:38:ILE:O	10:AJ:38:ILE:HG13	1.97	0.64
12:AL:6:THR:HG22	12:AL:9:GLN:HE21	1.61	0.64
13:AM:45:VAL:O	13:AM:48:LEU:HD22	1.97	0.64
16:AP:22:THR:HG22	16:AP:32:TYR:HA	1.78	0.64
17:AQ:29:HIS:HB3	17:AQ:33:GLY:N	2.13	0.64
19:AS:63:THR:CG2	19:AS:66:MET:HG2	2.22	0.64
20:AT:41:ILE:C	20:AT:43:LEU:H	2.00	0.64
31:B5:44:THR:HG21	48:BR:101:ALA:CB	2.27	0.64
35:BA:1894:C:O2'	35:BA:1895:C:H5'	1.98	0.64
35:BA:2225:A:H4'	35:BA:2226:C:H5'	1.78	0.64
35:BA:2881:C:H2'	35:BA:2882:A:H8	1.63	0.64
35:BA:510:C:H2'	35:BA:511:U:O4'	1.98	0.64
35:BA:878:A:H3'	35:BA:879:G:H8	1.63	0.64
36:BB:60:C:H2'	36:BB:61:G:H8	1.62	0.64
38:BD:35:LYS:HE2	38:BD:104:TYR:CB	2.27	0.64
38:BD:36:PRO:CG	38:BD:61:LEU:HD21	2.28	0.64
48:BR:9:LYS:NZ	48:BR:42:LYS:HB3	2.12	0.64
51:BU:61:TRP:HB3	51:BU:93:LYS:O	1.97	0.64
56:BZ:139:VAL:C	56:BZ:141:VAL:H	1.99	0.64
1:CA:1126:U:H2'	1:CA:1127:G:H8	1.62	0.64
1:CA:1466:C:H2'	1:CA:1467:G:O4'	1.98	0.64
3:CC:58:GLU:O	3:CC:64:VAL:HA	1.97	0.64
7:CG:101:LEU:O	7:CG:105:VAL:HG23	1.98	0.64
13:CM:45:VAL:O	13:CM:48:LEU:HD22	1.97	0.64
13:CM:9:ILE:N	13:CM:9:ILE:HD12	2.12	0.64
27:D1:76:ARG:HD3	27:D1:78:LYS:HZ3	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1280:G:C3'	35:DA:1281:G:H5''	2.28	0.64
35:DA:1387:C:H5'	35:DA:1469:A:H4'	1.79	0.64
35:DA:1697:G:H3'	35:DA:1698:A:C5'	2.27	0.64
36:DB:15:A:H3'	36:DB:16:G:C5'	2.27	0.64
42:DH:86:GLU:H	42:DH:86:GLU:CD	2.00	0.64
44:DN:1:MET:HG2	44:DN:2:LYS:N	2.12	0.64
36:DB:117:G:H5'	49:DS:55:ALA:HB1	1.80	0.64
50:DT:52:ILE:HG22	50:DT:61:PHE:CB	2.27	0.64
1:AA:1434:A:H2'	1:AA:1435:G:O4'	1.97	0.64
8:AH:17:THR:O	8:AH:19:VAL:N	2.31	0.64
9:AI:112:LYS:HA	9:AI:119:ALA:CB	2.27	0.64
12:AL:27:LEU:O	12:AL:29:GLY:N	2.30	0.64
27:B1:69:LYS:O	27:B1:73:LEU:HD12	1.98	0.64
30:B4:29:PRO:C	30:B4:31:ILE:H	2.00	0.64
34:B8:34:TRP:O	34:B8:35:GLN:HB2	1.97	0.64
35:BA:1643:G:H2'	35:BA:1644:C:H6	1.63	0.64
35:BA:1788:C:O2'	35:BA:1789:A:H5'	1.98	0.64
35:BA:1879:C:H2'	35:BA:1880:C:C5'	2.17	0.64
35:BA:2065:C:H1'	35:BA:2449:U:H3	1.62	0.64
35:BA:2726:U:H6	45:BO:67:LYS:HZ3	1.45	0.64
40:BF:7:TYR:HD2	40:BF:16:GLY:HA3	1.63	0.64
41:BG:114:ILE:HG22	41:BG:115:ARG:N	2.13	0.64
43:BI:2:LYS:HB2	43:BI:39:ALA:CB	2.26	0.64
46:BP:66:GLY:O	46:BP:68:GLN:N	2.30	0.64
49:BS:34:HIS:HE1	49:BS:55:ALA:HB2	1.63	0.64
50:BT:28:VAL:CG2	50:BT:47:GLY:N	2.61	0.64
50:BT:35:LYS:HE2	50:BT:41:ARG:HG3	1.79	0.64
35:BA:814:C:C5'	52:BV:86:GLY:HA3	2.28	0.64
56:BZ:136:PHE:C	56:BZ:136:PHE:CD1	2.71	0.64
1:CA:1242:C:H2'	1:CA:1243:C:C6	2.33	0.64
1:CA:1313:U:P	19:CS:6:LYS:HG3	2.37	0.64
1:CA:264:U:H2'	1:CA:265:G:O4'	1.98	0.64
1:CA:404:U:H2'	1:CA:405:U:C6	2.32	0.64
1:CA:514:C:H42	1:CA:537:G:H1	1.46	0.64
2:CB:19:HIS:N	2:CB:39:ILE:HG21	2.12	0.64
5:CE:42:GLY:HA3	5:CE:66:MET:HG2	1.80	0.64
10:CJ:38:ILE:HG13	10:CJ:38:ILE:O	1.96	0.64
25:CY:126:ARG:O	25:CY:129:ILE:N	2.31	0.64
27:D1:40:ARG:HG2	27:D1:41:ARG:H	1.61	0.64
27:D1:87:PRO:HB2	27:D1:91:LYS:HE3	1.78	0.64
33:D7:3:ARG:O	33:D7:6:GLN:NE2	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1406:U:H3'	35:DA:1407:C:H6	1.63	0.64
35:DA:2106:G:H2'	35:DA:2107:C:O4'	1.97	0.64
35:DA:2487:G:H2'	35:DA:2488:A:C8	2.33	0.64
35:DA:2673:G:H2'	35:DA:2674:G:H8	1.61	0.64
35:DA:2893:G:H5'	35:DA:2894:G:C5'	2.12	0.64
35:DA:646:A:H2'	35:DA:647:G:O4'	1.98	0.64
35:DA:807:U:O2'	35:DA:808:G:H5'	1.98	0.64
35:DA:838:C:H42	35:DA:940:G:H1	1.44	0.64
38:DD:226:MET:HB3	38:DD:230:ASP:CB	2.27	0.64
38:DD:183:ARG:HG3	38:DD:269:PHE:O	1.97	0.64
41:DG:19:LEU:O	41:DG:23:PHE:N	2.30	0.64
42:DH:87:LEU:N	42:DH:131:VAL:O	2.31	0.64
44:DN:17:ASP:O	44:DN:19:GLU:N	2.30	0.64
46:DP:126:VAL:HA	46:DP:145:PRO:HG2	1.78	0.64
49:DS:17:ARG:HG3	49:DS:18:ILE:HD12	1.80	0.64
1:AA:565:U:H3'	1:AA:566:G:H2'	1.80	0.64
1:AA:764:C:H2'	1:AA:765:G:H8	1.61	0.64
4:AD:15:GLU:HG2	4:AD:63:LYS:HG3	1.80	0.64
9:AI:111:ARG:O	9:AI:119:ALA:HB1	1.98	0.64
10:AJ:35:SER:O	10:AJ:72:VAL:HG13	1.97	0.64
11:AK:27:ASN:HA	11:AK:55:LYS:O	1.96	0.64
15:AO:15:PHE:HB2	15:AO:27:VAL:HG22	1.79	0.64
26:B0:25:ARG:HA	26:B0:29:GLN:NE2	2.12	0.64
33:B7:3:ARG:O	33:B7:6:GLN:NE2	2.31	0.64
35:BA:2206:G:C2	35:BA:2207:G:H5'	2.33	0.64
35:BA:2625:G:H2'	35:BA:2626:C:C6	2.33	0.64
35:BA:2795:G:N2	35:BA:2799:C:H5'	2.12	0.64
35:BA:313:C:H2'	35:BA:314:A:H8	1.63	0.64
35:BA:493:G:C3'	35:BA:494:G:H5''	2.27	0.64
37:BC:47:LEU:HA	37:BC:207:THR:HA	1.79	0.64
39:BE:36:ARG:HH22	39:BE:88:GLY:N	1.96	0.64
40:BF:3:GLU:HB2	40:BF:20:LEU:H	1.63	0.64
43:BI:14:ASP:O	43:BI:15:VAL:O	2.15	0.64
43:BI:68:LEU:HD23	43:BI:68:LEU:O	1.97	0.64
44:BN:9:VAL:CG1	44:BN:39:ARG:HH22	2.06	0.64
44:BN:46:VAL:HG11	44:BN:48:MET:HG3	1.78	0.64
46:BP:100:LEU:H	46:BP:100:LEU:HD22	1.63	0.64
48:BR:53:HIS:HA	48:BR:56:LYS:HB2	1.80	0.64
36:BB:49:C:OP1	49:BS:96:GLY:HA3	1.98	0.64
54:BX:54:VAL:C	54:BX:55:ASN:ND2	2.51	0.64
56:BZ:142:SER:N	56:BZ:144:LEU:HD23	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1003:G:H2'	1:CA:1004:A:O4'	1.97	0.64
1:CA:1202:G:H2'	1:CA:1203:C:H5'	1.80	0.64
1:CA:59:A:H5''	1:CA:60:A:H5''	1.77	0.64
1:CA:824:C:H4'	8:CH:1:MET:N	2.13	0.64
1:CA:954:G:H2'	1:CA:955:U:C6	2.32	0.64
2:CB:105:PHE:O	2:CB:108:ILE:N	2.31	0.64
2:CB:178:ARG:HB2	2:CB:178:ARG:HH11	1.63	0.64
8:CH:51:VAL:HG11	8:CH:60:ARG:HG2	1.80	0.64
9:CI:112:LYS:HA	9:CI:119:ALA:CB	2.28	0.64
3:CC:9:GLY:HA3	14:CN:49:HIS:HA	1.80	0.64
22:CV:38:U:H2'	22:CV:39:C:C6	2.33	0.64
35:DA:108:U:H2'	35:DA:109:G:C8	2.32	0.64
35:DA:2024:G:O2'	35:DA:2025:C:H5'	1.97	0.64
35:DA:286:C:H42	35:DA:355:G:H1	1.45	0.64
36:DB:61:G:H2'	36:DB:62:C:C6	2.33	0.64
40:DF:181:LEU:HG	40:DF:186:ILE:HD11	1.77	0.64
47:DQ:121:ALA:O	47:DQ:125:LEU:HD12	1.96	0.64
51:DU:13:LYS:O	51:DU:16:LYS:HB3	1.98	0.64
52:DV:52:VAL:C	52:DV:54:GLY:H	2.01	0.64
35:DA:2012:G:O2'	53:DW:96:ILE:HD11	1.97	0.64
54:DX:33:LYS:C	54:DX:35:THR:N	2.44	0.64
55:DY:37:VAL:HG13	55:DY:69:ALA:HB2	1.80	0.64
56:DZ:116:VAL:HG12	56:DZ:117:LEU:H	1.63	0.64
1:AA:1499:A:C2'	1:AA:1500:A:H5'	2.28	0.64
1:AA:601:C:H2'	1:AA:602:A:H8	1.62	0.64
1:AA:878:G:H5'	8:AH:89:PRO:CG	2.28	0.64
2:AB:102:LEU:N	2:AB:102:LEU:HD12	2.12	0.64
8:AH:89:PRO:HA	8:AH:92:ARG:NH1	2.13	0.64
9:AI:28:VAL:HG13	9:AI:63:ILE:O	1.98	0.64
15:AO:37:ASN:H	15:AO:37:ASN:ND2	1.95	0.64
23:AW:34:U:H1'	23:AW:36:A:N7	2.12	0.64
28:B2:12:GLU:O	28:B2:14:ARG:NH2	2.31	0.64
33:B7:34:ARG:O	33:B7:35:ARG:C	2.36	0.64
33:B7:5:TRP:NE1	33:B7:7:PRO:HG3	2.13	0.64
35:BA:1590:U:C3'	35:BA:1591:G:H5''	2.27	0.64
35:BA:718:A:H3'	35:BA:719:C:C6	2.32	0.64
40:BF:128:ALA:O	40:BF:142:TRP:NE1	2.31	0.64
41:BG:57:ALA:HB2	41:BG:90:LEU:HD23	1.80	0.64
44:BN:120:LEU:HD11	44:BN:122:VAL:CG2	2.26	0.64
46:BP:86:LYS:HD3	46:BP:117:GLU:HB2	1.79	0.64
50:BT:52:ILE:HG22	50:BT:61:PHE:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:48:ILE:C	50:BT:63:VAL:HG12	2.18	0.64
44:BN:38:HIS:O	51:BU:67:ALA:HB1	1.97	0.64
52:BV:3:ALA:CB	52:BV:14:VAL:HB	2.27	0.64
1:CA:1355:G:H2'	1:CA:1356:G:C8	2.32	0.64
1:CA:312:C:H2'	1:CA:313:A:C8	2.33	0.64
1:CA:547:A:H4'	1:CA:548:G:O5'	1.98	0.64
1:CA:571:U:H5''	1:CA:819:A:C2	2.32	0.64
1:CA:770:C:O2'	1:CA:771:G:H5'	1.97	0.64
2:CB:103:THR:HA	2:CB:180:LEU:HD11	1.79	0.64
2:CB:33:TYR:CD1	2:CB:43:ASP:HA	2.33	0.64
2:CB:75:LYS:CA	2:CB:78:GLN:HE21	2.01	0.64
4:CD:155:LEU:HB2	4:CD:158:ILE:HB	1.80	0.64
5:CE:126:ARG:C	5:CE:131:ILE:HD11	2.17	0.64
9:CI:63:ILE:HD12	9:CI:63:ILE:N	2.13	0.64
26:D0:25:ARG:HG2	26:D0:31:VAL:HG12	1.79	0.64
27:D1:83:GLU:CD	27:D1:86:SER:H	2.01	0.64
28:D2:53:LEU:HD12	35:DA:77:C:OP1	1.98	0.64
31:D5:40:LYS:HE2	31:D5:46:CYS:HB3	1.79	0.64
32:D6:12:GLU:CB	32:D6:23:THR:HG22	2.28	0.64
34:D8:50:LEU:N	34:D8:53:PRO:HG3	2.13	0.64
35:DA:1742:G:N7	35:DA:1743:C:N3	2.45	0.64
35:DA:2543:G:H2'	35:DA:2544:G:C8	2.32	0.64
35:DA:2801(A):A:C3'	35:DA:2802:G:H5'	2.28	0.64
38:DD:176:ARG:HH11	38:DD:176:ARG:HG2	1.63	0.64
35:DA:1567:A:H2'	38:DD:84:TYR:CE2	2.32	0.64
41:DG:105:LYS:HB2	41:DG:105:LYS:NZ	2.12	0.64
41:DG:94:LEU:O	41:DG:99:MET:HB2	1.98	0.64
45:DO:122:LEU:N	45:DO:122:LEU:HD12	2.13	0.64
48:DR:49:ASP:O	48:DR:50:HIS:C	2.36	0.64
49:DS:54:LEU:HD21	49:DS:59:LYS:O	1.98	0.64
56:DZ:115:GLY:HA3	56:DZ:175:VAL:O	1.98	0.64
1:AA:1091:U:H2'	1:AA:1093:A:OP2	1.97	0.64
1:AA:189(H):G:H2'	1:AA:189(I):G:H8	1.63	0.64
1:AA:264:U:H2'	1:AA:265:G:O4'	1.98	0.64
2:AB:111:ARG:NH1	2:AB:111:ARG:HG2	2.09	0.64
3:AC:112:SER:HB3	3:AC:115:LEU:HD12	1.79	0.64
1:AA:1190:G:H3'	3:AC:3:ASN:OD1	1.98	0.64
4:AD:120:LEU:H	4:AD:120:LEU:CD1	2.02	0.64
5:AE:103:GLY:O	5:AE:106:PRO:HD2	1.98	0.64
7:AG:16:LEU:HD13	9:AI:41:VAL:HG12	1.80	0.64
8:AH:10:LEU:N	8:AH:10:LEU:HD23	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AI:116:LYS:O	9:AI:118:LYS:N	2.31	0.64
12:AL:89:ARG:HH11	12:AL:89:ARG:C	2.01	0.64
13:AM:90:LEU:C	13:AM:92:HIS:N	2.51	0.64
16:AP:6:LEU:HB3	16:AP:17:TYR:CD2	2.32	0.64
35:BA:1257:C:O2'	40:BF:84:VAL:HG23	1.98	0.64
35:BA:237:C:H2'	35:BA:238:C:H6	1.63	0.64
35:BA:2872:G:C2	35:BA:2873:A:N6	2.66	0.64
35:BA:528:A:C2	35:BA:2042:A:H2'	2.33	0.64
35:BA:70:G:H2'	35:BA:113:G:O2'	1.98	0.64
35:BA:8:A:C4	35:BA:9:U:C5	2.86	0.64
36:BB:45:A:H1'	41:BG:95:ARG:NH2	2.12	0.64
38:BD:125:ILE:N	38:BD:125:ILE:HD12	2.12	0.64
40:BF:182:ASN:ND2	40:BF:185:ASP:OD2	2.31	0.64
40:BF:41:LEU:O	40:BF:44:ARG:HG3	1.98	0.64
42:BH:150:ALA:O	42:BH:152:ARG:N	2.31	0.64
46:BP:126:VAL:HA	46:BP:145:PRO:HB2	1.78	0.64
47:BQ:116:GLU:O	47:BQ:119:ARG:HB3	1.98	0.64
47:BQ:35:VAL:HG12	47:BQ:130:LYS:HB3	1.79	0.64
49:BS:62:LYS:N	49:BS:62:LYS:HD3	2.12	0.64
50:BT:91:ARG:O	50:BT:117:ASP:HB2	1.97	0.64
56:BZ:128:VAL:HG22	56:BZ:129:SER:H	1.63	0.64
1:CA:1499:A:H2'	1:CA:1500:A:H8	1.62	0.64
1:CA:542:G:H2'	1:CA:543:C:H6	1.61	0.64
2:CB:12:GLU:OE2	2:CB:214:ILE:HD11	1.98	0.64
4:CD:121:VAL:O	4:CD:134:ASP:HA	1.98	0.64
4:CD:15:GLU:HG2	4:CD:63:LYS:HG3	1.80	0.64
7:CG:146:GLU:HA	7:CG:149:ARG:HB2	1.80	0.64
9:CI:79:LEU:HD13	9:CI:79:LEU:O	1.98	0.64
10:CJ:39:PRO:HA	10:CJ:70:ARG:HA	1.79	0.64
19:CS:62:ILE:HD12	19:CS:66:MET:HG3	1.78	0.64
35:DA:1280:G:H3'	35:DA:1281:G:H5"	1.79	0.64
35:DA:1587:A:H2'	35:DA:1588:C:O4'	1.97	0.64
35:DA:2491:U:H4'	35:DA:2570:G:OP1	1.98	0.64
35:DA:519:U:H5"	53:DW:25:ARG:NH2	2.13	0.64
39:DE:36:ARG:HH22	39:DE:88:GLY:N	1.96	0.64
41:DG:37:VAL:HG11	41:DG:94:LEU:HD12	1.78	0.64
46:DP:81:GLN:NE2	46:DP:106:LEU:HA	2.12	0.64
46:DP:41:ARG:HD2	46:DP:41:ARG:N	2.12	0.64
50:DT:27:THR:C	50:DT:88:ILE:HD13	2.18	0.64
53:DW:12:ILE:CD1	53:DW:42:ARG:HH11	2.11	0.64
53:DW:51:LEU:C	53:DW:51:LEU:HD13	2.17	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1253:G:H5'	10:AJ:44:VAL:HG12	1.79	0.63
1:AA:1405:G:O2'	1:AA:1406:U:H5'	1.99	0.63
1:AA:560:U:O2'	1:AA:561:U:OP2	2.15	0.63
11:AK:58:PRO:HD3	11:AK:89:ALA:HB1	1.79	0.63
15:AO:36:ILE:HD12	15:AO:63:ARG:HD3	1.78	0.63
18:AR:36:ASN:ND2	18:AR:39:VAL:HB	2.14	0.63
20:AT:36:LEU:CD2	20:AT:36:LEU:H	2.11	0.63
20:AT:64:ASP:C	20:AT:66:ALA:H	2.02	0.63
26:B0:25:ARG:HG2	26:B0:31:VAL:HG12	1.79	0.63
28:B2:45:SER:HA	28:B2:48:HIS:HB2	1.80	0.63
34:B8:40:GLU:OE1	34:B8:44:LYS:HE3	1.98	0.63
35:BA:1301:A:HO2'	35:BA:1302:A:P	2.20	0.63
35:BA:2410:G:C2	35:BA:2411:A:H1'	2.33	0.63
34:B8:32:LEU:HD22	35:BA:2419:U:O5'	1.97	0.63
35:BA:2801(A):A:C4'	35:BA:2802:G:H2'	2.27	0.63
35:BA:2887:U:O2'	35:BA:2888:C:H5'	1.98	0.63
35:BA:556:G:H2'	35:BA:557:U:C6	2.33	0.63
43:BI:5:LEU:HD23	43:BI:5:LEU:N	2.13	0.63
45:BO:13:ASN:HD22	45:BO:97:ARG:CG	2.10	0.63
46:BP:80:TYR:CE1	46:BP:111:ARG:HB3	2.33	0.63
35:BA:587:C:C4	46:BP:33:ARG:HD2	2.33	0.63
52:BV:61:VAL:HG12	52:BV:62:LEU:H	1.62	0.63
54:BX:36:LYS:C	54:BX:38:GLU:N	2.51	0.63
56:BZ:7:ALA:H	56:BZ:62:PRO:HD3	1.63	0.63
1:CA:648:A:H2'	1:CA:649:G:H8	1.63	0.63
3:CC:119:ARG:HH21	3:CC:140:ARG:CZ	2.11	0.63
8:CH:17:THR:O	8:CH:19:VAL:N	2.31	0.63
9:CI:111:ARG:HG2	9:CI:112:LYS:N	2.11	0.63
27:D1:58:ILE:CD1	27:D1:59:THR:N	2.61	0.63
28:D2:21:LEU:HD22	28:D2:50:ILE:HG22	1.80	0.63
34:D8:62:LEU:HD13	35:DA:242:G:C5'	2.16	0.63
35:DA:2069:G:O2'	35:DA:2070:G:H5'	1.98	0.63
35:DA:2404:C:H2'	35:DA:2405:G:O4'	1.98	0.63
35:DA:2801(A):A:C4'	35:DA:2802:G:H2'	2.28	0.63
35:DA:2838:G:O2'	35:DA:2839:G:H5'	1.98	0.63
35:DA:577:G:H2'	35:DA:578:A:C8	2.33	0.63
38:DD:35:LYS:CE	38:DD:104:TYR:HB2	2.28	0.63
39:DE:116:VAL:O	39:DE:117:MET:CB	2.45	0.63
41:DG:36:LYS:H	41:DG:160:VAL:HB	1.62	0.63
41:DG:36:LYS:N	41:DG:160:VAL:HB	2.13	0.63
44:DN:67:LEU:HA	44:DN:88:GLU:HG3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:62:LEU:N	46:DP:62:LEU:HD13	2.12	0.63
35:DA:17:G:H4'	51:DU:25:TRP:CZ3	2.32	0.63
56:DZ:148:ASP:OD1	56:DZ:149:SER:N	2.29	0.63
5:AE:144:THR:O	5:AE:147:ASP:OD2	2.16	0.63
6:AF:49:ALA:HB1	18:AR:80:PRO:HG3	1.80	0.63
6:AF:61:LEU:O	6:AF:62:TRP:HB2	1.98	0.63
1:AA:1298:C:C4	7:AG:114:ARG:HD2	2.34	0.63
11:AK:19:ALA:HA	11:AK:32:ILE:HA	1.80	0.63
14:AN:23:ARG:HD3	14:AN:29:ARG:O	1.99	0.63
18:AR:44:LEU:O	18:AR:45:SER:HB3	1.98	0.63
26:B0:39:ARG:HH21	35:BA:2355:C:H1'	1.62	0.63
34:B8:50:LEU:N	34:B8:53:PRO:HG3	2.12	0.63
35:BA:1309:G:O2'	35:BA:1310:G:H5'	1.98	0.63
35:BA:582:G:H2'	35:BA:583:G:H8	1.62	0.63
35:BA:2051:A:H4'	39:BE:141:ILE:HD11	1.79	0.63
42:BH:149:ARG:HA	42:BH:162:ILE:CD1	2.28	0.63
46:BP:100:LEU:HD22	46:BP:100:LEU:N	2.13	0.63
40:BF:34:TRP:CB	46:BP:11:GLY:HA3	2.29	0.63
49:BS:90:GLY:HA2	49:BS:92:TYR:CD2	2.33	0.63
54:BX:59:VAL:C	54:BX:73:ARG:HA	2.18	0.63
54:BX:82:GLN:HG3	54:BX:83:VAL:N	2.13	0.63
54:BX:8:ILE:HD12	54:BX:8:ILE:N	2.13	0.63
1:CA:1325:C:O2	1:CA:1325:C:C2'	2.46	0.63
1:CA:184:G:C4'	1:CA:224:C:H4'	2.28	0.63
1:CA:956:U:O2'	1:CA:957:U:H5'	1.97	0.63
2:CB:111:ARG:HG2	2:CB:111:ARG:NH1	2.13	0.63
13:CM:65:LYS:C	13:CM:66:LEU:N	2.52	0.63
16:CP:2:VAL:HG23	16:CP:22:THR:O	1.99	0.63
28:D2:29:LYS:CA	28:D2:32:LEU:HB3	2.26	0.63
28:D2:41:ILE:C	28:D2:43:GLN:H	2.02	0.63
28:D2:49:LYS:CB	28:D2:53:LEU:HD22	2.18	0.63
35:DA:2030:A:H4'	35:DA:2031:A:H8	1.62	0.63
35:DA:2364:C:H2'	35:DA:2365:G:O4'	1.98	0.63
35:DA:250:G:H2'	35:DA:251:A:C8	2.34	0.63
35:DA:2722:G:H2'	35:DA:2723:C:C6	2.34	0.63
35:DA:2777:G:H5''	35:DA:2778:A:H5''	1.79	0.63
46:DP:122:PRO:HG3	46:DP:141:ALA:HB3	1.79	0.63
47:DQ:16:ARG:C	47:DQ:17:LEU:HD23	2.18	0.63
52:DV:43:GLU:HB2	52:DV:48:GLY:HA3	1.80	0.63
56:DZ:117:LEU:HA	56:DZ:173:ALA:O	1.98	0.63
56:DZ:58:VAL:HG13	56:DZ:67:LEU:C	2.19	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:73:GLN:CG	56:DZ:74:VAL:H	2.07	0.63
1:AA:1090:U:H2'	1:AA:1091:U:C6	2.33	0.63
1:AA:34:C:H2'	1:AA:35:G:C8	2.29	0.63
1:AA:758:G:H8	1:AA:758:G:O5'	1.81	0.63
2:AB:72:GLY:HA3	2:AB:165:VAL:CG1	2.28	0.63
2:AB:33:TYR:HB2	2:AB:41:ILE:HG22	1.80	0.63
3:AC:134:ILE:O	3:AC:137:ALA:HB3	1.98	0.63
4:AD:126:ILE:HG22	4:AD:127:THR:H	1.62	0.63
5:AE:7:GLU:O	5:AE:8:GLU:HB3	1.96	0.63
6:AF:37:VAL:HA	6:AF:65:VAL:HG12	1.80	0.63
11:AK:86:GLY:H	11:AK:112:THR:HG23	1.63	0.63
32:B6:11:LEU:CD1	32:B6:51:GLU:HB2	2.28	0.63
35:BA:1335:U:H2'	35:BA:1336:A:C8	2.33	0.63
35:BA:1771:C:H2'	35:BA:1772:G:H8	1.63	0.63
35:BA:1795:C:H42	35:BA:1824:G:H1	1.44	0.63
35:BA:1836:C:O2'	35:BA:1837:C:H5'	1.99	0.63
35:BA:286:C:H42	35:BA:355:G:H1	1.44	0.63
38:BD:226:MET:HB3	38:BD:230:ASP:CB	2.28	0.63
40:BF:148:LEU:HD21	40:BF:191:ARG:HD3	1.80	0.63
43:BI:130:TYR:HB2	43:BI:136:VAL:HG13	1.79	0.63
43:BI:83:ALA:HA	43:BI:89:TYR:HD1	1.64	0.63
56:BZ:127:LYS:H	56:BZ:164:ALA:CB	2.11	0.63
56:BZ:22:GLY:O	56:BZ:41:LEU:HG	1.97	0.63
1:CA:1007:C:H2'	1:CA:1008:C:C6	2.34	0.63
1:CA:1163:C:H2'	1:CA:1164:G:H8	1.63	0.63
1:CA:1436:U:H2'	1:CA:1437:C:H6	1.62	0.63
2:CB:144:ARG:O	2:CB:147:LYS:HB3	1.99	0.63
3:CC:43:LEU:CD2	3:CC:47:LEU:HD22	2.22	0.63
4:CD:187:ARG:HH11	4:CD:187:ARG:HG2	1.63	0.63
8:CH:26:VAL:HG22	8:CH:32:LYS:HZ2	1.61	0.63
8:CH:26:VAL:HG22	8:CH:32:LYS:HZ3	1.62	0.63
8:CH:97:VAL:HG13	8:CH:98:LYS:HG3	1.81	0.63
13:CM:97:PRO:HB3	13:CM:110:ARG:HD3	1.80	0.63
25:CY:21:LEU:HD21	25:CY:122:ALA:HA	1.81	0.63
35:DA:2020:A:O2'	35:DA:2021:C:H5'	1.98	0.63
35:DA:2197:U:H1'	35:DA:2198:A:C8	2.33	0.63
35:DA:2284:C:H2'	35:DA:2285:C:C5'	2.19	0.63
35:DA:2861:G:O2'	35:DA:2862:G:H5'	1.98	0.63
35:DA:809:G:O4'	35:DA:1254:A:H1'	1.99	0.63
38:DD:63:ARG:NH1	38:DD:86:PRO:HD2	2.14	0.63
41:DG:129:GLY:O	41:DG:131:TYR:N	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:56:ALA:HB1	41:DG:153:ARG:CD	2.28	0.63
35:DA:2820:A:C4'	48:DR:5:LYS:HE2	2.28	0.63
36:DB:7:G:H21	49:DS:38:GLN:NE2	1.97	0.63
49:DS:90:GLY:HA2	49:DS:92:TYR:CD2	2.32	0.63
54:DX:39:ILE:HD12	54:DX:40:LYS:N	2.13	0.63
54:DX:81:VAL:CG1	54:DX:85:PRO:HB2	2.28	0.63
1:AA:328:C:H2'	1:AA:328:C:O2	1.97	0.63
2:AB:22:LYS:HA	2:AB:22:LYS:HZ2	1.62	0.63
7:AG:91:VAL:HG13	7:AG:95:ARG:HD3	1.80	0.63
9:AI:50:LEU:HB3	9:AI:55:ALA:HB3	1.80	0.63
11:AK:103:LEU:N	11:AK:103:LEU:HD22	2.04	0.63
11:AK:88:GLY:O	11:AK:91:ARG:HB2	1.97	0.63
15:AO:18:PHE:CZ	15:AO:21:ASP:HB2	2.33	0.63
27:B1:87:PRO:HG2	27:B1:88:LYS:H	1.63	0.63
29:B3:4:LEU:HD21	29:B3:56:VAL:CG1	2.28	0.63
35:BA:1291:C:O2'	35:BA:1292:U:H5'	1.98	0.63
35:BA:2707:G:H5''	48:BR:68:ARG:HH21	1.63	0.63
35:BA:2801(A):A:C3'	35:BA:2802:G:H5'	2.29	0.63
35:BA:910:A:H62	47:BQ:12:GLN:HA	1.64	0.63
40:BF:199:TRP:O	40:BF:203:GLN:HG2	1.98	0.63
46:BP:16:ARG:NH1	46:BP:18:ARG:HG3	2.13	0.63
47:BQ:34:LEU:CD1	47:BQ:129:THR:HB	2.27	0.63
47:BQ:51:ARG:O	47:BQ:55:VAL:HG13	1.99	0.63
48:BR:29:LEU:HD23	48:BR:70:LEU:HD11	1.79	0.63
45:BO:119:PRO:HB2	50:BT:68:TYR:CD1	2.33	0.63
51:BU:30:LYS:HE3	51:BU:30:LYS:HA	1.81	0.63
51:BU:3:ARG:HH11	51:BU:3:ARG:CG	2.10	0.63
56:BZ:70:LEU:HB2	56:BZ:91:LEU:HD11	1.80	0.63
1:CA:1106:G:H2'	1:CA:1107:C:C6	2.33	0.63
1:CA:1518:A:C2	1:CA:1519:A:C2	2.86	0.63
1:CA:560:U:O2'	1:CA:561:U:OP2	2.13	0.63
2:CB:115:LEU:HG	2:CB:116:GLU:N	2.13	0.63
4:CD:19:LEU:HD23	4:CD:67:ILE:HA	1.80	0.63
4:CD:65:ARG:HH11	4:CD:72:GLU:N	1.97	0.63
1:CA:9:G:H5'	5:CE:122:GLU:OE2	1.98	0.63
12:CL:45:PRO:HB2	12:CL:49:ASN:O	1.98	0.63
12:CL:70:ILE:HG13	12:CL:100:ILE:HD12	1.80	0.63
13:CM:90:LEU:O	13:CM:92:HIS:N	2.28	0.63
17:CQ:77:VAL:O	17:CQ:78:GLU:HB3	1.98	0.63
20:CT:51:GLU:O	20:CT:55:ILE:HG12	1.99	0.63
25:CY:83:ILE:O	25:CY:86:SER:N	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D2:32:LEU:HG	28:D2:33:MET:N	2.13	0.63
35:DA:1031:G:N2	35:DA:1124:C:H1'	2.13	0.63
35:DA:1590:U:C3'	35:DA:1591:G:H5''	2.28	0.63
35:DA:2056:G:N2	35:DA:2057:A:H1'	2.14	0.63
35:DA:2199:A:H5'	35:DA:2200:C:OP2	1.99	0.63
35:DA:2745:C:H2'	35:DA:2746:U:C6	2.34	0.63
35:DA:2772:C:H2'	35:DA:2773:C:H6	1.62	0.63
35:DA:807:U:H2'	35:DA:808:G:C8	2.33	0.63
38:DD:175:LEU:HD23	38:DD:176:ARG:N	2.13	0.63
38:DD:183:ARG:HG2	38:DD:183:ARG:HH11	1.63	0.63
39:DE:154:LYS:HA	39:DE:154:LYS:HE3	1.79	0.63
40:DF:7:TYR:HD2	40:DF:16:GLY:HA3	1.63	0.63
41:DG:144:ILE:CG1	41:DG:145:THR:N	2.56	0.63
42:DH:136:ILE:O	42:DH:137:ASP:HB2	1.97	0.63
44:DN:57:ALA:HB1	44:DN:60:ILE:HD11	1.81	0.63
46:DP:86:LYS:HD3	46:DP:117:GLU:HB2	1.78	0.63
49:DS:34:HIS:HE1	49:DS:55:ALA:HB2	1.63	0.63
35:DA:1754:C:H5'	50:DT:101:PHE:CE1	2.33	0.63
51:DU:18:LEU:CD2	51:DU:22:LYS:HE2	2.28	0.63
51:DU:6:THR:O	51:DU:8:VAL:N	2.28	0.63
52:DV:37:VAL:HG12	52:DV:38:LEU:N	2.13	0.63
35:DA:814:C:C5'	52:DV:86:GLY:HA3	2.27	0.63
55:DY:31:LEU:HD23	55:DY:36:ALA:HB3	1.80	0.63
1:AA:32:A:H2'	1:AA:33:A:H8	1.59	0.63
1:AA:375:U:H2'	1:AA:376:G:C8	2.33	0.63
10:AJ:32:ALA:HB1	10:AJ:75:ILE:CG1	2.28	0.63
23:AW:49:C:O2'	23:AW:60:A:H4'	1.98	0.63
35:BA:1150:C:O2'	35:BA:1151:G:H5'	1.98	0.63
35:BA:116:C:O2'	35:BA:117:G:H5'	1.99	0.63
35:BA:1327:C:H2'	35:BA:1328:G:O4'	1.99	0.63
35:BA:2312:U:C2'	35:BA:2313:C:H5''	2.27	0.63
35:BA:2626:C:H2'	35:BA:2627:G:C8	2.34	0.63
35:BA:575:A:H2'	35:BA:576:U:H5'	1.80	0.63
35:BA:743:G:O2'	35:BA:744:G:H5'	1.97	0.63
38:BD:45:ASN:CG	38:BD:46:GLN:N	2.51	0.63
42:BH:31:GLY:O	42:BH:79:VAL:HG11	1.97	0.63
45:BO:104:ARG:HH21	50:BT:33:LYS:HE3	1.63	0.63
45:BO:47:ILE:HG23	45:BO:48:PRO:HD2	1.80	0.63
50:BT:99:LEU:O	50:BT:99:LEU:HD13	1.97	0.63
52:BV:71:LEU:HD13	52:BV:72:VAL:N	2.13	0.63
1:CA:1348:U:H4'	9:CI:120:ARG:HD2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:155:C:H2'	1:CA:156:G:C8	2.34	0.63
1:CA:328:C:H2'	1:CA:328:C:O2	1.98	0.63
1:CA:404:U:H2'	1:CA:405:U:H6	1.64	0.63
2:CB:165:VAL:CG2	2:CB:166:ASP:N	2.59	0.63
3:CC:173:VAL:N	3:CC:174:PRO:HD3	2.14	0.63
3:CC:73:PRO:HA	3:CC:76:VAL:CG1	2.28	0.63
9:CI:3:GLN:HG2	9:CI:20:ARG:HH21	1.64	0.63
24:CX:13:A:O2'	24:CX:14:U:H5'	1.98	0.63
35:DA:1518:U:H2'	35:DA:1519:G:O4'	1.98	0.63
35:DA:1523:U:H2'	35:DA:1524:G:H8	1.63	0.63
35:DA:759:G:H2'	35:DA:760:G:C8	2.33	0.63
36:DB:52:A:HO2'	36:DB:53:A:H8	1.46	0.63
39:DE:197:ILE:HD11	39:DE:199:ARG:HH22	1.62	0.63
40:DF:3:GLU:HB2	40:DF:20:LEU:H	1.64	0.63
43:DI:2:LYS:HB2	43:DI:39:ALA:CB	2.26	0.63
45:DO:111:PHE:HB3	45:DO:114:ILE:CD1	2.29	0.63
46:DP:111:ARG:HA	46:DP:128:HIS:ND1	2.13	0.63
1:AA:1341:U:O2'	1:AA:1342:C:H5'	1.98	0.63
1:AA:648:A:H2'	1:AA:649:G:H8	1.63	0.63
1:AA:950:U:H2'	1:AA:951:G:C8	2.34	0.63
2:AB:172:ILE:H	2:AB:172:ILE:CD1	2.07	0.63
3:AC:16:ARG:CB	3:AC:16:ARG:HH11	2.12	0.63
4:AD:127:THR:HG22	4:AD:149:ALA:H	1.63	0.63
7:AG:121:ALA:HA	7:AG:124:LEU:HD12	1.81	0.63
1:AA:668:G:O4'	15:AO:49:ASP:HB2	1.97	0.63
27:B1:53:VAL:HG12	27:B1:58:ILE:HG22	1.80	0.63
32:B6:12:GLU:CB	32:B6:23:THR:HG22	2.29	0.63
35:BA:1605:C:H5'	35:BA:1610:A:N6	2.14	0.63
35:BA:1991:U:H2'	35:BA:1992:G:H5'	1.79	0.63
35:BA:45:C:OP2	35:BA:215:G:H2'	1.97	0.63
27:B1:33:LYS:HB2	35:BA:2395:C:O2'	1.99	0.63
41:BG:67:LYS:H	41:BG:67:LYS:CD	2.09	0.63
42:BH:162:ILE:HD12	42:BH:162:ILE:C	2.18	0.63
46:BP:95:VAL:HG23	46:BP:125:VAL:HB	1.79	0.63
46:BP:130:PHE:CD2	46:BP:130:PHE:N	2.67	0.63
50:BT:96:ARG:HH11	50:BT:96:ARG:CG	2.12	0.63
1:CA:1470:G:O2'	1:CA:1471:G:H5'	1.99	0.63
1:CA:570:G:H2'	1:CA:571:U:C6	2.33	0.63
4:CD:13:ARG:HD3	4:CD:38:TYR:O	1.98	0.63
5:CE:76:ILE:HD11	5:CE:142:LEU:HD11	1.80	0.63
6:CF:98:LEU:H	6:CF:98:LEU:HD12	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CK:29:ILE:C	11:CK:29:ILE:HD12	2.18	0.63
6:CF:49:ALA:HB1	18:CR:80:PRO:HG3	1.79	0.63
25:CY:69:GLN:HA	25:CY:97:ASP:O	1.98	0.63
35:DA:1153:C:H2'	35:DA:1154:G:O4'	1.98	0.63
35:DA:1887:C:H3'	35:DA:1888:G:H5''	1.80	0.63
35:DA:1758:G:N7	35:DA:2695:C:H4'	2.13	0.63
35:DA:271(V):G:H2'	35:DA:271(W):G:O4'	1.97	0.63
40:DF:117:ARG:HH21	40:DF:187:VAL:HA	1.64	0.63
35:DA:1131:G:OP1	44:DN:80:GLY:HA2	1.99	0.63
54:DX:36:LYS:C	54:DX:38:GLU:N	2.51	0.63
54:DX:77:LYS:HE3	54:DX:78:LYS:N	2.12	0.63
47:DQ:140:ALA:HB1	56:DZ:99:TYR:H	1.64	0.63
2:AB:115:LEU:HG	2:AB:116:GLU:N	2.12	0.63
22:AV:38:U:H2'	22:AV:39:C:H6	1.62	0.63
25:AY:176:ALA:O	25:AY:180:GLU:HG3	1.99	0.63
31:B5:20:ARG:HH12	53:BW:15:ARG:CZ	2.12	0.63
35:BA:1049:C:H2'	35:BA:1050:A:H8	1.63	0.63
35:BA:1196:C:H2'	35:BA:1197:G:C8	2.33	0.63
35:BA:1938:A:H2	35:BA:2590:A:H1'	1.62	0.63
35:BA:2203:U:O4'	38:BD:151:LYS:HE3	1.99	0.63
35:BA:2467:C:H2'	35:BA:2468:G:H5'	1.79	0.63
35:BA:534:U:O3'	51:BU:46:ALA:HB2	1.99	0.63
35:BA:581:C:O2'	35:BA:582:G:H5'	1.99	0.63
41:BG:76:SER:CB	41:BG:84:LYS:H	2.12	0.63
41:BG:42:GLY:HA2	41:BG:89:GLY:HA2	1.79	0.63
44:BN:17:ASP:C	44:BN:19:GLU:N	2.52	0.63
45:BO:3:GLN:HB2	45:BO:4:PRO:HD2	1.79	0.63
46:BP:85:LEU:HA	46:BP:88:LEU:CB	2.26	0.63
47:BQ:82:ARG:NH1	47:BQ:82:ARG:HG2	2.13	0.63
35:BA:975(A):G:OP1	52:BV:79:VAL:HG13	1.97	0.63
54:BX:57:LEU:HB2	54:BX:76:ARG:CD	2.29	0.63
55:BY:37:VAL:HG13	55:BY:69:ALA:HB2	1.81	0.63
56:BZ:60:GLU:O	56:BZ:62:PRO:HD3	1.99	0.63
1:CA:342:C:O2'	1:CA:343:U:H5'	1.98	0.63
6:CF:68:PRO:HG3	6:CF:71:ARG:NH2	2.11	0.63
7:CG:16:LEU:HD13	9:CI:41:VAL:HG12	1.81	0.63
10:CJ:8:LEU:HA	10:CJ:96:ILE:HG22	1.79	0.63
17:CQ:45:HIS:HB2	17:CQ:69:LYS:HE2	1.78	0.63
18:CR:78:LEU:O	18:CR:79:LEU:HG	1.98	0.63
19:CS:36:ARG:HH12	19:CS:75:ALA:HB3	1.64	0.63
20:CT:37:SER:O	20:CT:40:ALA:HB3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:D7:30:VAL:HG13	33:D7:33:ARG:HH22	1.64	0.63
35:DA:2061:G:H5''	35:DA:2503:A:C2	2.33	0.63
35:DA:2103:C:H2'	35:DA:2104:G:H5''	1.81	0.63
35:DA:2521:C:N4	35:DA:2544:G:H1	1.97	0.63
35:DA:426:C:O2'	35:DA:427:U:H5'	1.99	0.63
36:DB:74:U:H2'	36:DB:75:G:C5'	2.18	0.63
38:DD:80:ALA:HB3	38:DD:94:LEU:CD1	2.28	0.63
40:DF:110:LEU:HD21	40:DF:181:LEU:HD23	1.80	0.63
44:DN:43:THR:O	44:DN:46:VAL:N	2.31	0.63
46:DP:80:TYR:CD1	46:DP:111:ARG:HB3	2.33	0.63
47:DQ:34:LEU:HD11	47:DQ:129:THR:CB	2.25	0.63
53:DW:75:TYR:N	53:DW:75:TYR:CD1	2.63	0.63
1:AA:1030(A):G:H1'	1:AA:1031:G:H1	1.63	0.63
1:AA:334:C:O2'	1:AA:335:C:H5'	1.98	0.63
2:AB:194:PRO:O	2:AB:196:LEU:N	2.29	0.63
16:AP:6:LEU:HD12	16:AP:6:LEU:N	2.14	0.63
6:AF:62:TRP:CB	18:AR:35:ARG:HH12	2.12	0.63
20:AT:37:SER:O	20:AT:40:ALA:HB3	1.99	0.63
23:AW:11:A:H2'	23:AW:12:G:C8	2.33	0.63
27:B1:25:LYS:CB	27:B1:37:ILE:HD11	2.28	0.63
31:B5:49:CYS:HB2	31:B5:59:GLU:OE1	1.98	0.63
35:BA:1453:U:H5'	48:BR:63:ARG:NE	2.14	0.63
35:BA:1946:U:O2'	35:BA:1947:C:H5'	1.98	0.63
35:BA:1998:G:H2'	35:BA:1999:C:C6	2.33	0.63
35:BA:189:G:H2'	35:BA:205:G:H22	1.62	0.63
35:BA:543:C:N4	35:BA:551:G:N1	2.47	0.63
35:BA:646:A:H2'	35:BA:647:G:O4'	1.98	0.63
39:BE:48:GLN:HG2	39:BE:78:LEU:HD12	1.81	0.63
39:BE:93:VAL:C	39:BE:95:ILE:H	2.02	0.63
41:BG:19:LEU:HD21	41:BG:175:LEU:HD13	1.79	0.63
42:BH:86:GLU:H	42:BH:86:GLU:CD	2.02	0.63
44:BN:58:ASP:O	44:BN:60:ILE:N	2.30	0.63
50:BT:94:ALA:HB1	50:BT:99:LEU:HD23	1.81	0.63
35:BA:2012:G:O2'	53:BW:96:ILE:HD11	1.99	0.63
54:BX:72:LYS:CE	54:BX:74:PRO:HB3	2.23	0.63
56:BZ:150:LEU:HD23	56:BZ:171:ILE:HG13	1.80	0.63
56:BZ:4:ARG:HG2	56:BZ:58:VAL:O	1.98	0.63
56:BZ:3:TYR:C	56:BZ:57:ILE:HG23	2.19	0.63
1:CA:17:U:H2'	1:CA:18:C:H6	1.64	0.63
1:CA:245:C:O2'	1:CA:246:A:H5'	1.99	0.63
1:CA:786:G:H1	1:CA:796:C:N4	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:115:LEU:HB2	2:CB:145:LEU:HD11	1.79	0.63
3:CC:119:ARG:NH2	3:CC:140:ARG:CZ	2.62	0.63
5:CE:43:LEU:HD12	5:CE:44:GLY:N	2.14	0.63
8:CH:88:LYS:HB3	8:CH:89:PRO:HD2	1.81	0.63
10:CJ:32:ALA:H	10:CJ:78:ASN:CG	2.01	0.63
15:CO:37:ASN:H	15:CO:37:ASN:ND2	1.95	0.63
16:CP:82:GLN:NE2	16:CP:82:GLN:N	2.45	0.63
17:CQ:60:ILE:HG12	17:CQ:61:GLU:O	1.98	0.63
27:D1:88:LYS:C	27:D1:90:ILE:N	2.46	0.63
28:D2:20:GLU:O	28:D2:23:LYS:N	2.31	0.63
28:D2:26:ARG:HD3	54:DX:5:TYR:CD1	2.34	0.63
28:D2:52:ASP:O	28:D2:54:LYS:N	2.31	0.63
34:D8:38:GLY:O	34:D8:39:LYS:HB3	1.97	0.63
35:DA:1300:U:H1'	35:DA:1626:G:C2	2.34	0.63
35:DA:2387:U:H5'	35:DA:2388:A:OP2	1.98	0.63
35:DA:2469:A:H2	35:DA:2481:G:H21	1.46	0.63
35:DA:2814:C:H2'	35:DA:2815:C:H6	1.61	0.63
35:DA:492:A:H2'	35:DA:493:G:O4'	1.98	0.63
35:DA:565:C:O3'	52:DV:81:TYR:HE1	1.80	0.63
35:DA:863:A:O2'	35:DA:864:G:H5'	1.98	0.63
36:DB:65:C:C2'	36:DB:66:A:H5'	2.28	0.63
38:DD:117:VAL:HG22	38:DD:118:VAL:H	1.61	0.63
39:DE:36:ARG:HH22	39:DE:88:GLY:H	1.45	0.63
46:DP:130:PHE:N	46:DP:130:PHE:CD2	2.67	0.63
48:DR:74:LYS:O	48:DR:77:ARG:N	2.32	0.63
48:DR:9:LYS:HE3	48:DR:43:GLU:OE2	1.98	0.63
52:DV:14:VAL:HG12	52:DV:15:GLU:H	1.64	0.63
31:D5:25:LEU:CD1	53:DW:19:LEU:HB3	2.28	0.63
54:DX:54:VAL:C	54:DX:55:ASN:ND2	2.52	0.63
1:AA:1149:C:H2'	1:AA:1150:U:H6	1.63	0.63
1:AA:237:C:H4'	17:AQ:25:ARG:NH1	2.13	0.63
3:AC:119:ARG:HH21	3:AC:140:ARG:CZ	2.12	0.63
9:AI:71:SER:HA	9:AI:74:ILE:HD12	1.81	0.63
10:AJ:5:ARG:HG3	10:AJ:73:ASP:OD1	1.99	0.63
13:AM:90:LEU:O	13:AM:92:HIS:N	2.29	0.63
25:AY:73:GLN:HG3	25:AY:74:ASN:N	2.14	0.63
27:B1:26:ARG:CB	27:B1:34:THR:OG1	2.46	0.63
32:B6:32:ASN:ND2	32:B6:33:LYS:H	1.95	0.63
34:B8:32:LEU:O	34:B8:34:TRP:N	2.29	0.63
35:BA:2197:U:H1'	35:BA:2198:A:C8	2.33	0.63
39:BE:119:ARG:HD2	39:BE:120:TRP:NE1	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:114:VAL:CG2	40:BF:115:ALA:H	2.11	0.63
46:BP:126:VAL:HA	46:BP:145:PRO:CG	2.29	0.63
48:BR:20:LEU:HD12	48:BR:21:TYR:N	2.14	0.63
54:BX:68:ARG:HG3	54:BX:69:TYR:CD1	2.34	0.63
55:BY:86:ARG:NH2	55:BY:95:LYS:HZ2	1.97	0.63
56:BZ:58:VAL:HG13	56:BZ:68:PRO:N	2.14	0.63
1:CA:101:A:H2'	1:CA:102:G:H8	1.63	0.63
1:CA:1162:C:H2'	1:CA:1163:C:C6	2.33	0.63
1:CA:123:C:OP1	1:CA:312:C:H5'	1.99	0.63
1:CA:865:A:H5'	1:CA:1078:U:H5	1.64	0.63
1:CA:880:C:O2'	1:CA:881:G:H5'	1.99	0.63
1:CA:903:G:H2'	1:CA:904:C:H6	1.62	0.63
2:CB:100:GLY:HA2	2:CB:103:THR:HB	1.80	0.63
4:CD:117:ALA:O	4:CD:121:VAL:HG23	1.99	0.63
4:CD:127:THR:CG2	4:CD:149:ALA:H	2.12	0.63
4:CD:56:VAL:O	4:CD:58:LEU:N	2.32	0.63
5:CE:102:ALA:HB1	5:CE:106:PRO:CG	2.29	0.63
7:CG:36:LYS:HA	7:CG:39:ALA:HB2	1.80	0.63
7:CG:42:ILE:HA	7:CG:45:ASP:HB2	1.79	0.63
8:CH:1:MET:HE2	8:CH:2:LEU:N	2.13	0.63
9:CI:126:SER:O	9:CI:127:LYS:HB3	1.99	0.63
35:DA:1150:C:O2'	35:DA:1151:G:H5'	1.99	0.63
35:DA:1625:C:H2'	35:DA:1626:G:O4'	1.99	0.63
35:DA:2065:C:H1'	35:DA:2449:U:H3	1.63	0.63
35:DA:1938:A:H2	35:DA:2590:A:H1'	1.64	0.63
35:DA:2609:U:H4'	35:DA:2609:U:OP1	1.97	0.63
35:DA:811:U:H1'	35:DA:1251:C:C5'	2.29	0.63
38:DD:155:LEU:O	38:DD:156:ALA:C	2.37	0.63
43:DI:140:LEU:HD12	43:DI:141:LYS:H	1.60	0.63
43:DI:96:ASP:HA	43:DI:99:GLU:HB3	1.81	0.63
46:DP:124:LYS:HA	46:DP:143:GLY:N	2.14	0.63
46:DP:99:LEU:HD12	46:DP:102:ARG:HD2	1.81	0.63
49:DS:88:ASP:OD2	49:DS:89:ARG:N	2.29	0.63
51:DU:61:TRP:HB3	51:DU:93:LYS:O	1.98	0.63
54:DX:55:ASN:HD22	54:DX:55:ASN:N	1.97	0.63
54:DX:81:VAL:HG12	54:DX:82:GLN:O	1.99	0.63
54:DX:81:VAL:HG13	54:DX:85:PRO:HB2	1.80	0.63
54:DX:83:VAL:O	54:DX:84:ALA:HB3	1.98	0.63
56:DZ:104:PHE:CE2	56:DZ:122:ARG:HA	2.34	0.63
56:DZ:28:MET:HE1	56:DZ:59:LEU:HD12	1.81	0.63
56:DZ:53:ILE:HG22	56:DZ:71:VAL:HB	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:16:A:N1	1:AA:919:A:H2	1.96	0.62
1:AA:20:U:O2'	1:AA:21:G:H5'	1.99	0.62
1:AA:386:C:C2'	1:AA:387:U:H5'	2.29	0.62
1:AA:580:U:H2'	1:AA:581:G:O4'	1.99	0.62
1:AA:924:C:H2'	1:AA:925:G:H8	1.64	0.62
3:AC:111:LEU:CD2	3:AC:146:ALA:HB2	2.29	0.62
4:AD:79:PHE:HA	4:AD:93:PHE:CD2	2.34	0.62
9:AI:17:VAL:HG22	9:AI:63:ILE:HG23	1.81	0.62
12:AL:9:GLN:O	12:AL:11:VAL:N	2.32	0.62
1:AA:1227:A:OP2	13:AM:111:LYS:HE3	1.98	0.62
18:AR:22:VAL:HA	18:AR:25:THR:OG1	1.99	0.62
20:AT:100:ILE:HG22	20:AT:102:GLY:H	1.62	0.62
20:AT:23:ARG:O	20:AT:27:LYS:HB2	1.98	0.62
23:AW:10:G:O5'	23:AW:10:G:H8	1.81	0.62
34:B8:46:ARG:O	34:B8:47:LYS:HB3	1.97	0.62
35:BA:1159:U:H2'	35:BA:1160:G:H5'	1.79	0.62
35:BA:1434:A:H2'	35:BA:1435:G:C8	2.33	0.62
35:BA:1665:A:O2'	35:BA:1666:G:H5'	1.99	0.62
35:BA:18:C:H2'	35:BA:19:C:C6	2.32	0.62
35:BA:2777:G:H5''	35:BA:2778:A:H5''	1.81	0.62
35:BA:598:G:H5'	46:BP:15:ARG:HB2	1.80	0.62
36:BB:35:U:O2'	36:BB:36:C:H5'	1.99	0.62
38:BD:186:HIS:CD2	38:BD:188:GLU:H	2.14	0.62
39:BE:35:GLN:HE22	39:BE:37:ARG:NH2	1.97	0.62
43:BI:64:GLU:OE1	43:BI:67:ARG:HB2	1.99	0.62
44:BN:9:VAL:HG12	44:BN:10:GLU:N	2.14	0.62
46:BP:16:ARG:CZ	46:BP:18:ARG:HB2	2.29	0.62
35:BA:196:A:H5''	46:BP:46:LYS:NZ	2.13	0.62
46:BP:66:GLY:O	46:BP:68:GLN:HG2	1.98	0.62
50:BT:10:VAL:C	50:BT:12:SER:N	2.46	0.62
52:BV:61:VAL:CG2	52:BV:100:ARG:HG2	2.29	0.62
54:BX:51:VAL:CG1	54:BX:80:ILE:H	2.12	0.62
56:BZ:149:SER:HB3	56:BZ:173:ALA:CA	2.26	0.62
56:BZ:126:VAL:CA	56:BZ:164:ALA:HB3	2.29	0.62
56:BZ:58:VAL:HA	56:BZ:68:PRO:HA	1.81	0.62
1:CA:104:G:O2'	1:CA:105:G:H5'	1.99	0.62
1:CA:519:C:H2'	1:CA:520:A:H8	1.64	0.62
1:CA:661:G:H2'	1:CA:662:G:H8	1.64	0.62
1:CA:77:G:H1	1:CA:92:C:H42	1.47	0.62
1:CA:943:U:O5'	1:CA:943:U:H6	1.82	0.62
2:CB:73:THR:HG22	2:CB:94:ASN:HA	1.79	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:119:LYS:HB2	12:CL:120:TYR:HD1	1.63	0.62
25:CY:169:ILE:O	25:CY:172:ALA:CB	2.47	0.62
26:D0:40:GLN:HE21	26:D0:43:THR:HA	1.64	0.62
27:D1:76:ARG:O	27:D1:77:ALA:HB3	1.99	0.62
35:DA:1428:C:N4	35:DA:1569:A:H3'	2.14	0.62
35:DA:2111:C:H5'	35:DA:2112:G:OP1	1.98	0.62
35:DA:230:U:H2'	35:DA:230:U:O2	1.97	0.62
35:DA:285:C:C2'	35:DA:286:C:H5''	2.27	0.62
35:DA:692:C:O2'	35:DA:693:C:H5'	1.98	0.62
38:DD:10:THR:HG23	38:DD:13:ARG:CB	2.28	0.62
38:DD:120:GLY:O	38:DD:131:LEU:HB3	1.99	0.62
38:DD:227:ASN:HB3	38:DD:228:PRO:HD2	1.81	0.62
40:DF:182:ASN:ND2	40:DF:185:ASP:OD2	2.32	0.62
40:DF:52:LYS:HD3	40:DF:57:VAL:HA	1.80	0.62
41:DG:109:VAL:HG12	41:DG:140:ILE:O	1.98	0.62
45:DO:2:ILE:HD12	45:DO:6:THR:HG21	1.79	0.62
45:DO:77:ILE:HD12	50:DT:73:GLU:O	1.99	0.62
47:DQ:60:ARG:C	47:DQ:60:ARG:HD3	2.19	0.62
47:DQ:71:ASP:O	47:DQ:73:PRO:HD3	1.99	0.62
35:DA:1275:A:C5	48:DR:16:HIS:ND1	2.68	0.62
48:DR:9:LYS:NZ	48:DR:42:LYS:HB3	2.13	0.62
48:DR:2:ARG:CZ	48:DR:5:LYS:HE3	2.28	0.62
49:DS:38:GLN:OE1	49:DS:47:THR:HG23	1.98	0.62
50:DT:48:ILE:HD12	50:DT:48:ILE:N	2.13	0.62
52:DV:71:LEU:HD13	52:DV:72:VAL:N	2.13	0.62
1:AA:1457:G:H2'	1:AA:1458:G:C8	2.35	0.62
1:AA:514:C:H42	1:AA:537:G:H1	1.45	0.62
1:AA:77:G:H1	1:AA:92:C:H42	1.46	0.62
1:AA:93:G:O2'	1:AA:96:U:H5'	1.99	0.62
6:AF:78:GLU:O	6:AF:81:ILE:HG13	1.99	0.62
12:AL:45:PRO:HB2	12:AL:49:ASN:O	1.99	0.62
16:AP:68:ASP:O	16:AP:71:ARG:HB3	1.99	0.62
29:B3:1:MET:SD	29:B3:38:GLU:HG2	2.40	0.62
33:B7:15:THR:HG22	33:B7:16:HIS:CD2	2.34	0.62
34:B8:38:GLY:O	34:B8:39:LYS:HB3	1.98	0.62
35:BA:1300:U:H1'	35:BA:1626:G:C2	2.34	0.62
35:BA:1324:G:H3'	35:BA:1325:G:C5'	2.30	0.62
35:BA:1809:A:H2'	35:BA:1810:A:C8	2.33	0.62
35:BA:2715:C:H2'	35:BA:2716:U:H6	1.64	0.62
36:BB:82:G:H2'	36:BB:83:G:H8	1.64	0.62
39:BE:2:LYS:HE2	39:BE:95:ILE:HG22	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BH:70:THR:CG2	42:BH:74:ASN:HD21	2.10	0.62
45:BO:77:ILE:HD12	50:BT:73:GLU:O	1.99	0.62
50:BT:91:ARG:O	50:BT:93:ARG:N	2.32	0.62
51:BU:92:ARG:C	51:BU:94:ASN:H	2.01	0.62
54:BX:65:ARG:NE	54:BX:66:LEU:H	1.97	0.62
56:BZ:129:SER:HB3	56:BZ:132:ASN:ND2	2.14	0.62
1:CA:1040:U:H2'	1:CA:1041:A:C8	2.35	0.62
1:CA:1205:U:O2'	3:CC:195:VAL:HG23	1.99	0.62
1:CA:312:C:H2'	1:CA:313:A:H8	1.63	0.62
1:CA:586:C:O2'	1:CA:587:G:H5'	1.99	0.62
1:CA:1190:G:H8	3:CC:3:ASN:HD21	1.45	0.62
9:CI:17:VAL:HG22	9:CI:63:ILE:HG23	1.80	0.62
17:CQ:95:TYR:O	17:CQ:97:SER:N	2.31	0.62
19:CS:63:THR:HG22	19:CS:66:MET:CG	2.26	0.62
31:D5:46:CYS:HB3	31:D5:48:GLU:OE2	1.99	0.62
34:D8:32:LEU:HD22	35:DA:2419:U:O5'	1.98	0.62
34:D8:49:VAL:HB	34:D8:53:PRO:HD3	1.80	0.62
35:DA:313:C:H2'	35:DA:314:A:H8	1.64	0.62
35:DA:834:C:H2'	35:DA:835:A:H8	1.64	0.62
39:DE:197:ILE:CD1	39:DE:199:ARG:HH22	2.12	0.62
41:DG:11:TYR:O	41:DG:13:GLU:N	2.33	0.62
41:DG:15:VAL:HG13	41:DG:175:LEU:CB	2.27	0.62
41:DG:96:ARG:HA	41:DG:99:MET:CE	2.29	0.62
43:DI:99:GLU:O	43:DI:103:ARG:HB2	1.99	0.62
48:DR:12:ARG:O	48:DR:13:HIS:HB3	1.99	0.62
48:DR:60:LEU:HD23	48:DR:61:HIS:N	2.14	0.62
45:DO:104:ARG:HH21	50:DT:33:LYS:HE3	1.63	0.62
55:DY:13:VAL:HG21	55:DY:28:LYS:HG2	1.80	0.62
56:DZ:108:PRO:HB3	56:DZ:144:LEU:H	1.62	0.62
1:AA:1126:U:H2'	1:AA:1127:G:H8	1.63	0.62
1:AA:1325:C:O2	1:AA:1325:C:C2'	2.47	0.62
1:AA:1329:A:OP1	13:AM:28:ALA:HB3	1.98	0.62
1:AA:597:G:C2'	1:AA:598:U:H5'	2.29	0.62
1:AA:783:C:C5	1:AA:784:C:H5	2.16	0.62
8:AH:26:VAL:HG22	8:AH:32:LYS:HZ3	1.64	0.62
9:AI:10:ARG:NH2	9:AI:107:ARG:HD3	2.14	0.62
1:AA:1342:C:H1'	9:AI:124:GLN:NE2	2.14	0.62
10:AJ:36:GLY:O	10:AJ:72:VAL:HG22	1.99	0.62
15:AO:53:HIS:CE1	15:AO:57:LEU:HD21	2.35	0.62
16:AP:49:LEU:HD12	16:AP:50:LYS:H	1.64	0.62
29:B3:45:GLY:HA3	35:BA:851:U:O2'	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:29:LYS:O	33:B7:32:LYS:HB3	1.99	0.62
35:BA:1428:C:N4	35:BA:1569:A:H3'	2.13	0.62
35:BA:2030:A:H4'	35:BA:2031:A:H8	1.63	0.62
35:BA:94(A):G:H2'	35:BA:95:G:C5'	2.26	0.62
38:BD:4:LYS:HZ1	38:BD:20:ASP:HA	1.62	0.62
44:BN:57:ALA:O	44:BN:58:ASP:O	2.17	0.62
44:BN:67:LEU:HA	44:BN:88:GLU:HG3	1.80	0.62
45:BO:46:ALA:H	45:BO:54:GLU:HG2	1.64	0.62
46:BP:111:ARG:HG3	46:BP:128:HIS:ND1	2.14	0.62
47:BQ:9:TYR:O	47:BQ:10:ARG:CG	2.43	0.62
49:BS:38:GLN:OE1	49:BS:47:THR:HG23	1.99	0.62
50:BT:28:VAL:HG21	50:BT:46:GLU:HA	1.81	0.62
51:BU:96:ALA:O	51:BU:98:LEU:N	2.32	0.62
54:BX:82:GLN:CD	54:BX:83:VAL:H	2.01	0.62
55:BY:38:ILE:HG22	55:BY:39:VAL:N	2.14	0.62
56:BZ:28:MET:SD	56:BZ:37:VAL:HG11	2.39	0.62
1:CA:1511:G:C6	1:CA:1512:U:N3	2.68	0.62
1:CA:222:U:H2'	1:CA:223:U:C6	2.34	0.62
1:CA:357:G:OP1	1:CA:367:U:H5''	1.99	0.62
1:CA:834:C:H2'	1:CA:835:U:C6	2.34	0.62
3:CC:134:ILE:HD11	3:CC:153:VAL:HG21	1.80	0.62
7:CG:11:GLN:HE21	7:CG:12:LEU:H	1.47	0.62
7:CG:91:VAL:HG13	7:CG:95:ARG:HD3	1.80	0.62
9:CI:111:ARG:O	9:CI:119:ALA:HB1	2.00	0.62
20:CT:23:ARG:O	20:CT:27:LYS:HB2	1.99	0.62
32:D6:11:LEU:CD1	32:D6:51:GLU:HB2	2.28	0.62
35:DA:1192:G:O2'	35:DA:1193:G:H5'	1.98	0.62
35:DA:1677:A:H2'	35:DA:1678:G:H8	1.61	0.62
35:DA:1760:A:O2'	35:DA:1761:C:H5'	1.99	0.62
35:DA:2775:A:O2'	35:DA:2776:A:H5'	1.98	0.62
35:DA:510:C:H2'	35:DA:511:U:O4'	2.00	0.62
35:DA:543:C:N4	35:DA:551:G:N1	2.47	0.62
35:DA:863:A:C2'	35:DA:864:G:H5'	2.28	0.62
37:DC:58:VAL:HG21	37:DC:166:ASP:N	2.14	0.62
38:DD:161:THR:O	38:DD:196:VAL:HG23	1.99	0.62
44:DN:56:ASN:HA	44:DN:125:GLY:H	1.64	0.62
44:DN:46:VAL:HG22	44:DN:47:ALA:N	2.14	0.62
46:DP:126:VAL:HA	46:DP:145:PRO:CB	2.28	0.62
46:DP:16:ARG:NH1	46:DP:18:ARG:HG3	2.15	0.62
35:DA:631:A:O2'	46:DP:67:MET:HB3	1.99	0.62
52:DV:3:ALA:CB	52:DV:14:VAL:HB	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:DW:95:ILE:O	53:DW:95:ILE:HG13	1.99	0.62
1:AA:1355:G:H2'	1:AA:1356:G:C8	2.34	0.62
1:AA:404:U:H2'	1:AA:405:U:C6	2.34	0.62
1:AA:880:C:OP2	12:AL:6:THR:HG21	1.99	0.62
5:AE:147:ASP:HA	5:AE:150:ARG:HD2	1.79	0.62
6:AF:98:LEU:H	6:AF:98:LEU:HD12	1.63	0.62
7:AG:11:GLN:NE2	7:AG:12:LEU:H	1.96	0.62
16:AP:14:ASN:N	16:AP:15:PRO:CD	2.62	0.62
18:AR:53:ARG:NH2	18:AR:60:ALA:N	2.45	0.62
25:AY:108:GLU:HA	25:AY:111:ARG:HG3	1.81	0.62
25:AY:64:ARG:HH21	25:AY:103:ILE:HD11	1.64	0.62
35:BA:863:A:O2'	35:BA:864:G:H5'	1.98	0.62
35:BA:881:G:N2	35:BA:896:A:H62	1.96	0.62
39:BE:134:ILE:N	39:BE:134:ILE:HD13	2.14	0.62
39:BE:120:TRP:CE2	39:BE:155:LYS:HB3	2.34	0.62
41:BG:41:GLN:C	41:BG:43:LEU:H	2.02	0.62
44:BN:58:ASP:O	44:BN:60:ILE:HG13	1.99	0.62
50:BT:96:ARG:HG3	50:BT:98:LYS:O	1.99	0.62
52:BV:89:GLN:HE21	52:BV:89:GLN:HA	1.65	0.62
53:BW:12:ILE:CD1	53:BW:42:ARG:HH11	2.13	0.62
55:BY:28:LYS:CA	55:BY:39:VAL:H	2.12	0.62
55:BY:42:VAL:HB	55:BY:65:ALA:HB3	1.81	0.62
1:CA:1438:G:H2'	1:CA:1439:C:H6	1.62	0.62
1:CA:503:C:H2'	1:CA:504:C:H6	1.62	0.62
1:CA:702:A:H3'	1:CA:703:G:H5'	1.81	0.62
7:CG:93:PRO:HA	7:CG:96:GLN:NE2	2.14	0.62
8:CH:11:THR:HA	8:CH:14:ARG:HH12	1.63	0.62
35:DA:1424:G:H2'	35:DA:1425:G:O4'	1.99	0.62
35:DA:2710:C:OP1	48:DR:15:SER:HB2	2.00	0.62
35:DA:2866:U:C5	35:DA:2868:A:H1'	2.34	0.62
35:DA:64:A:H2'	35:DA:65:C:H6	1.64	0.62
38:DD:206:LEU:HA	38:DD:211:ARG:NH1	2.14	0.62
39:DE:5:LEU:HD22	39:DE:197:ILE:HG22	1.81	0.62
41:DG:16:ARG:HB3	41:DG:17:PRO:HD3	1.81	0.62
46:DP:126:VAL:HA	46:DP:145:PRO:CG	2.28	0.62
35:DA:2413:G:H21	46:DP:70:GLN:HE21	1.46	0.62
1:AA:1168:A:H2'	1:AA:1169:A:C8	2.35	0.62
1:AA:1246:C:H2'	1:AA:1247:U:C6	2.35	0.62
1:AA:222:U:H2'	1:AA:223:U:C6	2.34	0.62
1:AA:473:G:H5''	16:AP:81:ARG:HE	1.65	0.62
1:AA:67:C:H2'	1:AA:68:G:H8	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:132:LYS:O	2:AB:135:GLN:HB2	2.00	0.62
2:AB:144:ARG:O	2:AB:147:LYS:HB3	1.98	0.62
2:AB:168:THR:HG21	2:AB:191:ASP:OD1	1.99	0.62
2:AB:28:PHE:CE1	2:AB:31:TYR:HB2	2.35	0.62
2:AB:19:HIS:N	2:AB:39:ILE:HG21	2.15	0.62
4:AD:145:GLU:HG2	4:AD:184:LYS:CG	2.26	0.62
6:AF:21:LEU:HA	6:AF:24:GLU:HG2	1.82	0.62
8:AH:36:LEU:O	8:AH:39:LEU:HB3	1.98	0.62
9:AI:114:TYR:CE1	10:AJ:59:SER:HA	2.34	0.62
10:AJ:32:ALA:H	10:AJ:78:ASN:CG	2.02	0.62
12:AL:83:VAL:CG2	12:AL:84:LEU:H	2.09	0.62
15:AO:11:VAL:HG21	15:AO:34:LEU:HD23	1.80	0.62
18:AR:38:GLU:O	18:AR:41:LYS:HB3	2.00	0.62
28:B2:41:ILE:O	28:B2:42:GLY:C	2.38	0.62
35:BA:1350:C:O2'	35:BA:1351:C:H5'	1.99	0.62
35:BA:1760:A:O2'	35:BA:1761:C:H5'	1.99	0.62
35:BA:2562:U:C2'	35:BA:2563:U:H5'	2.28	0.62
38:BD:264:LYS:HG3	38:BD:265:PRO:HD2	1.81	0.62
39:BE:115:GLY:HA2	39:BE:157:ALA:HB1	1.82	0.62
40:BF:88:VAL:HG22	40:BF:89:VAL:N	2.13	0.62
41:BG:41:GLN:NE2	41:BG:153:ARG:HB3	2.14	0.62
42:BH:41:MET:CE	42:BH:55:PRO:HD2	2.29	0.62
44:BN:32:THR:HG22	44:BN:37:LYS:HB3	1.82	0.62
46:BP:45:LEU:HD23	46:BP:46:LYS:N	2.14	0.62
47:BQ:29:PHE:HB2	47:BQ:65:PHE:CE2	2.34	0.62
48:BR:45:ARG:HG3	48:BR:46:GLY:N	2.10	0.62
50:BT:27:THR:OG1	50:BT:87:ASP:HA	2.00	0.62
56:BZ:107:THR:HG21	56:BZ:111:VAL:HG11	1.82	0.62
1:CA:1329:A:OP1	13:CM:28:ALA:HB3	1.98	0.62
1:CA:1443:G:H8	1:CA:1443:G:OP2	1.83	0.62
2:CB:24:TRP:CG	2:CB:25:ASN:N	2.67	0.62
3:CC:53:ALA:HB2	3:CC:115:LEU:CD2	2.30	0.62
8:CH:119:LEU:N	8:CH:119:LEU:CD2	2.63	0.62
8:CH:39:LEU:O	8:CH:44:PHE:HB2	2.00	0.62
9:CI:4:TYR:HB2	9:CI:19:LEU:HD12	1.81	0.62
12:CL:25:PRO:HD2	12:CL:98:TYR:OH	1.99	0.62
27:D1:76:ARG:HD3	27:D1:78:LYS:NZ	2.14	0.62
35:DA:1350:C:O2'	35:DA:1351:C:H5'	2.00	0.62
35:DA:2712:U:H1'	35:DA:2712(A):A:H8	1.64	0.62
35:DA:380:U:H2'	35:DA:381:G:H8	1.65	0.62
35:DA:910:A:H62	47:DQ:12:GLN:HA	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:209:ALA:C	38:DD:210:GLY:O	2.35	0.62
38:DD:94:LEU:CD1	38:DD:94:LEU:H	2.06	0.62
41:DG:142:PRO:HG2	41:DG:143:GLU:H	1.64	0.62
45:DO:26:LYS:HB2	45:DO:30:ALA:HB2	1.81	0.62
1:AA:1040:U:H2'	1:AA:1041:A:C8	2.35	0.62
1:AA:702:A:H3'	1:AA:703:G:H5'	1.81	0.62
2:AB:51:LEU:HB3	2:AB:55:PHE:HE2	1.65	0.62
4:AD:104:VAL:HG21	4:AD:140:VAL:HG21	1.80	0.62
7:AG:36:LYS:HA	7:AG:39:ALA:HB2	1.81	0.62
9:AI:105:ASP:HB3	9:AI:107:ARG:HG3	1.82	0.62
11:AK:87:THR:HA	11:AK:91:ARG:HG3	1.81	0.62
12:AL:24:VAL:HG13	12:AL:98:TYR:HE2	1.65	0.62
1:AA:472:A:H1'	16:AP:82:GLN:OE1	2.00	0.62
18:AR:58:LEU:HD12	18:AR:58:LEU:N	2.08	0.62
19:AS:6:LYS:CD	19:AS:7:LYS:HD2	2.30	0.62
35:BA:1190:G:O5'	46:BP:35:HIS:HA	2.00	0.62
35:BA:1289:C:H2'	35:BA:1290:C:C6	2.35	0.62
35:BA:1344:G:H4'	35:BA:1384:A:C5	2.34	0.62
35:BA:2019:A:H4'	51:BU:34:LYS:HD2	1.82	0.62
35:BA:26:G:H1'	35:BA:515:A:N6	2.12	0.62
35:BA:2759:G:O2'	35:BA:2760:C:H5'	1.99	0.62
35:BA:971:C:H2'	35:BA:972:G:O4'	2.00	0.62
44:BN:10:GLU:CG	44:BN:11:PRO:HD2	2.30	0.62
44:BN:62:VAL:HG22	44:BN:66:LYS:HG3	1.80	0.62
45:BO:105:GLU:HA	45:BO:108:GLU:CG	2.29	0.62
46:BP:112:LEU:C	46:BP:112:LEU:HD13	2.20	0.62
51:BU:64:ARG:NH2	51:BU:64:ARG:CB	2.61	0.62
55:BY:36:ALA:HB1	55:BY:67:LEU:O	1.99	0.62
1:CA:1195:C:H5'	1:CA:1196:U:OP2	1.99	0.62
1:CA:447:G:N2	1:CA:488:C:H42	1.98	0.62
1:CA:556:C:O2'	1:CA:557:G:H5'	1.99	0.62
7:CG:135:VAL:O	7:CG:138:LYS:HB3	1.99	0.62
7:CG:150:ALA:C	7:CG:152:ALA:H	2.02	0.62
9:CI:50:LEU:HB3	9:CI:55:ALA:HB3	1.81	0.62
25:CY:29:ARG:HB2	25:CY:32:ARG:NH2	2.15	0.62
25:CY:3:LEU:N	25:CY:3:LEU:CD1	2.62	0.62
35:DA:1190:G:H4'	46:DP:35:HIS:CB	2.29	0.62
35:DA:1515:G:C2'	35:DA:1516:C:H5'	2.29	0.62
25:CY:130:ARG:NH2	35:DA:1943:U:OP1	2.31	0.62
32:D6:39:TYR:CE1	35:DA:2347:C:H4'	2.31	0.62
35:DA:2707:G:H2'	35:DA:2708:G:C8	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:585:G:H2'	35:DA:1251:C:N4	2.13	0.62
35:DA:632:A:H2'	35:DA:633:A:C8	2.34	0.62
35:DA:709:U:H2'	35:DA:710:G:C8	2.35	0.62
36:DB:82:G:H2'	36:DB:83:G:H8	1.64	0.62
38:DD:57:GLY:HA2	38:DD:214:TRP:O	1.99	0.62
40:DF:28:ILE:N	40:DF:28:ILE:HD13	2.15	0.62
41:DG:107:LEU:HA	41:DG:111:LEU:HD12	1.81	0.62
41:DG:29:TRP:C	41:DG:31:VAL:N	2.50	0.62
42:DH:86:GLU:HA	42:DH:132:ARG:HB3	1.82	0.62
43:DI:130:TYR:HB2	43:DI:136:VAL:HG13	1.82	0.62
46:DP:66:GLY:O	46:DP:68:GLN:N	2.33	0.62
46:DP:86:LYS:HB3	46:DP:117:GLU:C	2.20	0.62
1:AA:1007:C:H2'	1:AA:1008:C:C6	2.35	0.62
1:AA:1064:G:H1'	1:AA:1065:U:H5''	1.82	0.62
1:AA:1348:U:H4'	9:AI:120:ARG:HD2	1.82	0.62
1:AA:174:C:O2'	1:AA:175:C:H5'	2.00	0.62
1:AA:710:G:H5''	6:AF:54:LYS:HE3	1.81	0.62
3:AC:182:ILE:HG23	3:AC:203:PHE:HA	1.82	0.62
1:AA:796:C:P	11:AK:123:LYS:HZ2	2.23	0.62
17:AQ:77:VAL:O	17:AQ:78:GLU:HB3	1.98	0.62
25:AY:122:ALA:O	25:AY:126:ARG:HG3	2.00	0.62
25:AY:147:LEU:HD23	25:AY:148:HIS:N	2.14	0.62
35:BA:1018:C:O2'	35:BA:1019:U:H5'	2.00	0.62
35:BA:1330:C:O2'	35:BA:1331:A:H5'	2.00	0.62
35:BA:2334:G:H5'	49:BS:13:ARG:CG	2.25	0.62
35:BA:271(J):C:H2'	35:BA:271(K):U:H5''	1.82	0.62
35:BA:2864:G:H5'	35:BA:2864:G:H8	1.63	0.62
35:BA:805:G:H22	35:BA:828:U:H5''	1.64	0.62
38:BD:206:LEU:HA	38:BD:211:ARG:NH1	2.14	0.62
42:BH:137:ASP:O	42:BH:138:LYS:HB2	2.00	0.62
43:BI:4:ILE:HA	43:BI:17:GLN:O	2.00	0.62
44:BN:110:GLY:HA2	44:BN:114:ARG:NH2	2.15	0.62
44:BN:129:PRO:O	44:BN:130:HIS:HB3	2.00	0.62
35:BA:993:G:H5'	52:BV:75:PHE:CZ	2.35	0.62
56:BZ:9:TYR:HE2	56:BZ:35:ARG:HD2	1.65	0.62
1:CA:1206:G:O4'	3:CC:194:GLY:HA2	2.00	0.62
1:CA:189(H):G:H2'	1:CA:189(I):G:H8	1.63	0.62
1:CA:386:C:C2'	1:CA:387:U:H5'	2.30	0.62
4:CD:126:ILE:HG22	4:CD:127:THR:N	2.14	0.62
7:CG:113:GLU:CB	7:CG:119:ARG:HG2	2.30	0.62
8:CH:18:ARG:H	8:CH:78:GLN:NE2	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CJ:4:ILE:CB	10:CJ:74:ILE:HD11	2.28	0.62
11:CK:99:GLN:HE22	11:CK:105:VAL:HG11	1.65	0.62
15:CO:28:GLN:O	15:CO:32:LEU:HG	1.99	0.62
17:CQ:76:LEU:HD12	17:CQ:77:VAL:H	1.65	0.62
19:CS:53:ASN:ND2	19:CS:55:LYS:H	1.98	0.62
23:CW:3:C:C2'	23:CW:4:G:H5''	2.23	0.62
25:CY:132:ILE:O	25:CY:133:ARG:C	2.38	0.62
25:CY:59:THR:O	25:CY:66:LEU:HD12	1.99	0.62
27:D1:85:LEU:CB	27:D1:87:PRO:HD3	2.29	0.62
29:D3:45:GLY:HA3	35:DA:851:U:O2'	1.99	0.62
35:DA:2292:C:O2'	35:DA:2293:C:H5'	1.99	0.62
35:DA:2544:G:H2'	35:DA:2545:G:H8	1.63	0.62
35:DA:2866:U:C6	35:DA:2868:A:H1'	2.35	0.62
35:DA:523:C:O2'	35:DA:524:U:H5'	2.00	0.62
35:DA:755:C:H2'	35:DA:756:C:H6	1.65	0.62
35:DA:796:C:H2'	35:DA:797:C:H6	1.64	0.62
35:DA:881:G:N2	35:DA:896:A:H62	1.96	0.62
35:DA:8:A:C4	35:DA:9:U:C5	2.86	0.62
36:DB:91:C:H2'	36:DB:92:C:C6	2.34	0.62
39:DE:24:THR:CG2	39:DE:184:VAL:HG23	2.27	0.62
41:DG:111:LEU:HA	41:DG:114:ILE:CD1	2.29	0.62
45:DO:120:GLU:HG3	45:DO:122:LEU:HD11	1.81	0.62
46:DP:16:ARG:CZ	46:DP:18:ARG:HB2	2.30	0.62
46:DP:70:GLN:HG3	46:DP:71:VAL:H	1.65	0.62
48:DR:37:THR:CG2	48:DR:40:LYS:HE2	2.30	0.62
51:DU:102:GLU:O	51:DU:105:VAL:HG23	2.00	0.62
51:DU:64:ARG:NH2	51:DU:64:ARG:CB	2.58	0.62
52:DV:19:LYS:HB3	52:DV:96:ILE:O	2.00	0.62
53:DW:11:ARG:HH22	53:DW:98:LYS:HB3	1.64	0.62
55:DY:36:ALA:HB1	55:DY:67:LEU:O	2.00	0.62
1:AA:1495:U:O2'	1:AA:1496:C:H5'	2.00	0.62
1:AA:376:G:OP2	16:AP:67:THR:HG21	1.99	0.62
1:AA:460:G:H21	1:AA:472:A:H62	1.48	0.62
1:AA:542:G:H2'	1:AA:543:C:H6	1.62	0.62
2:AB:72:GLY:HA3	2:AB:165:VAL:HG22	1.82	0.62
3:AC:43:LEU:CD2	3:AC:47:LEU:HD22	2.21	0.62
4:AD:127:THR:CG2	4:AD:149:ALA:H	2.12	0.62
13:AM:19:LEU:CA	13:AM:22:ILE:HD13	2.28	0.62
27:B1:20:ARG:HH22	27:B1:41:ARG:HE	1.46	0.62
34:B8:8:LYS:HB3	34:B8:12:LYS:HE3	1.81	0.62
35:BA:1344:G:H4'	35:BA:1384:A:N7	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2103:C:H3'	35:BA:2104:G:H5''	1.82	0.62
35:BA:437:G:O2'	35:BA:438:G:H5'	1.98	0.62
35:BA:492:A:H2'	35:BA:493:G:O4'	2.00	0.62
40:BF:161:GLU:O	40:BF:164:ARG:HB2	1.99	0.62
40:BF:181:LEU:HG	40:BF:186:ILE:HD11	1.81	0.62
30:B4:14:ILE:CB	41:BG:5:VAL:HG13	2.30	0.62
49:BS:24:LEU:HB2	49:BS:85:VAL:HB	1.80	0.62
54:BX:83:VAL:O	54:BX:84:ALA:HB3	2.00	0.62
55:BY:90:LEU:HG	55:BY:91:GLU:HG2	1.80	0.62
56:BZ:125:LEU:C	56:BZ:126:VAL:HG22	2.20	0.62
1:CA:403:C:O2'	1:CA:404:U:H5'	1.99	0.62
2:CB:168:THR:HG23	2:CB:192:SER:OG	1.99	0.62
4:CD:12:CYS:SG	4:CD:19:LEU:O	2.57	0.62
5:CE:147:ASP:HA	5:CE:150:ARG:HD2	1.81	0.62
11:CK:44:SER:O	11:CK:47:VAL:HB	1.99	0.62
23:CW:43:G:H2'	23:CW:44:A:H5''	1.82	0.62
25:CY:123:GLU:O	25:CY:124:GLU:C	2.38	0.62
25:CY:55:ILE:HD12	25:CY:55:ILE:N	2.15	0.62
35:DA:1703:G:H2'	35:DA:1704:G:C8	2.34	0.62
35:DA:2040:C:H2'	35:DA:2041:U:C6	2.34	0.62
35:DA:2368:C:H2'	35:DA:2369:A:H8	1.65	0.62
35:DA:685:A:C4	35:DA:689:A:N6	2.67	0.62
35:DA:736:C:H2'	35:DA:737:C:H6	1.65	0.62
36:DB:35:U:O2'	36:DB:36:C:H5'	1.99	0.62
38:DD:229:VAL:HG23	38:DD:230:ASP:N	2.15	0.62
39:DE:142:GLY:C	39:DE:143:ASN:ND2	2.53	0.62
40:DF:81:PRO:HG2	40:DF:82:ILE:H	1.65	0.62
41:DG:7:LEU:CA	41:DG:10:LYS:HB2	2.25	0.62
41:DG:25:TYR:CD1	41:DG:30:GLU:HG3	2.35	0.62
43:DI:88:ILE:HG22	43:DI:89:TYR:N	2.13	0.62
44:DN:107:LEU:HD12	44:DN:108:PRO:O	1.99	0.62
54:DX:89:ILE:HG22	54:DX:89:ILE:O	1.99	0.62
55:DY:31:LEU:CB	55:DY:36:ALA:H	2.13	0.62
1:AA:1090:U:H2'	1:AA:1091:U:H6	1.65	0.62
1:AA:1106:G:H2'	1:AA:1107:C:H6	1.65	0.62
1:AA:1411:C:H2'	1:AA:1412:C:C6	2.35	0.62
1:AA:148:G:H1	1:AA:174:C:H42	1.46	0.62
1:AA:294:U:H2'	1:AA:295:C:C6	2.34	0.62
2:AB:102:LEU:O	2:AB:105:PHE:HB2	2.00	0.62
3:AC:92:ALA:N	3:AC:99:VAL:HG21	2.14	0.62
4:AD:155:LEU:HB2	4:AD:158:ILE:HB	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:176:LEU:HG	4:AD:177:ASP:N	2.12	0.62
4:AD:19:LEU:HD23	4:AD:67:ILE:HA	1.82	0.62
6:AF:28:ARG:HH11	6:AF:28:ARG:HG3	1.65	0.62
7:AG:146:GLU:HA	7:AG:149:ARG:HB2	1.80	0.62
13:AM:91:ARG:HH11	19:AS:81:ARG:NH2	1.87	0.62
25:AY:179:LYS:O	25:AY:183:ILE:HG13	2.00	0.62
25:AY:84:ARG:C	25:AY:86:SER:H	2.03	0.62
34:B8:59:LYS:C	34:B8:61:LEU:N	2.53	0.62
35:BA:1518:U:H2'	35:BA:1519:G:O4'	2.00	0.62
35:BA:1567:A:H2'	38:BD:84:TYR:CE2	2.29	0.62
35:BA:2078:C:H2'	35:BA:2079:U:C6	2.35	0.62
35:BA:268:C:N4	35:BA:424:G:H1	1.97	0.62
35:BA:2830:G:C5'	39:BE:58:ARG:HH22	2.12	0.62
35:BA:541:C:H2'	35:BA:542:C:C5	2.34	0.62
35:BA:64:A:H2'	35:BA:65:C:C6	2.35	0.62
35:BA:755:C:H2'	35:BA:756:C:H6	1.65	0.62
35:BA:74:A:H4'	35:BA:75:G:O5'	2.00	0.62
38:BD:143:HIS:CE1	38:BD:192:THR:HG23	2.35	0.62
40:BF:96:ASP:OD1	40:BF:98:SER:HB3	1.99	0.62
43:BI:129:THR:OG1	43:BI:135:GLU:HB3	2.00	0.62
35:BA:1246:A:OP2	46:BP:16:ARG:NH2	2.33	0.62
47:BQ:103:MET:CE	47:BQ:125:LEU:HD21	2.29	0.62
35:BA:1278:A:H5''	48:BR:36:THR:HG22	1.81	0.62
51:BU:13:LYS:O	51:BU:16:LYS:HB3	1.99	0.62
51:BU:44:ASN:O	51:BU:47:TYR:HB3	2.00	0.62
1:CA:1064:G:H1'	1:CA:1065:U:H5''	1.81	0.62
1:CA:1226:C:H2'	13:CM:103:THR:OG1	2.00	0.62
1:CA:1267:C:H2'	1:CA:1267:C:O2	2.00	0.62
1:CA:1281:U:H4'	1:CA:1282:C:OP2	2.00	0.62
1:CA:280:C:C2	17:CQ:38:ARG:HG3	2.35	0.62
1:CA:652:U:H1'	1:CA:653:A:H2	1.64	0.62
1:CA:783:C:C5	1:CA:784:C:H5	2.17	0.62
1:CA:950:U:H2'	1:CA:951:G:C8	2.34	0.62
2:CB:69:LEU:HB2	2:CB:159:PRO:HG2	1.82	0.62
2:CB:72:GLY:CA	2:CB:165:VAL:HG22	2.30	0.62
4:CD:9:CYS:SG	4:CD:31:CYS:O	2.58	0.62
7:CG:50:ILE:HG22	7:CG:56:GLN:O	1.98	0.62
8:CH:122:ARG:HA	8:CH:125:ARG:CB	2.28	0.62
9:CI:48:GLU:N	9:CI:49:PRO:HD2	2.15	0.62
1:CA:963:G:N2	10:CJ:55:LYS:HZ3	1.96	0.62
10:CJ:94:VAL:HG12	10:CJ:95:GLU:N	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1049:C:H2'	35:DA:1050:A:H8	1.64	0.62
35:DA:1114:G:H2'	35:DA:1115:G:C5'	2.30	0.62
35:DA:1790:C:H2'	35:DA:1791:A:C8	2.35	0.62
35:DA:2103:C:H3'	35:DA:2104:G:H5''	1.81	0.62
35:DA:2175:C:H2'	35:DA:2176:A:C5'	2.29	0.62
35:DA:2241:A:H2'	35:DA:2242:G:C8	2.34	0.62
35:DA:2570:G:H2'	35:DA:2571:C:H6	1.63	0.62
35:DA:598:G:H5'	46:DP:15:ARG:HB2	1.80	0.62
35:DA:648:G:H2'	35:DA:649:G:C8	2.34	0.62
35:DA:841:A:H2'	35:DA:842:G:C8	2.35	0.62
38:DD:36:PRO:CG	38:DD:61:LEU:HD21	2.27	0.62
39:DE:38:THR:C	39:DE:40:GLU:H	2.03	0.62
41:DG:76:SER:HB3	41:DG:83:ARG:HB3	1.81	0.62
42:DH:67:LEU:O	42:DH:71:LEU:HD22	2.00	0.62
50:DT:23:ARG:HG2	50:DT:120:ARG:HH12	1.64	0.62
51:DU:58:ARG:HA	51:DU:61:TRP:CE3	2.35	0.62
56:DZ:108:PRO:HB3	56:DZ:142:SER:O	2.00	0.62
56:DZ:76:LEU:HA	56:DZ:84:GLU:H	1.65	0.62
1:AA:110:C:H2'	1:AA:111:G:O4'	2.00	0.62
2:AB:115:LEU:HB2	2:AB:145:LEU:HD11	1.82	0.62
2:AB:54:THR:O	2:AB:58:ILE:HG13	1.99	0.62
3:AC:173:VAL:N	3:AC:174:PRO:HD3	2.15	0.62
3:AC:70:VAL:HG12	3:AC:72:LYS:H	1.63	0.62
8:AH:119:LEU:H	8:AH:119:LEU:HD23	1.65	0.62
8:AH:18:ARG:H	8:AH:78:GLN:NE2	1.97	0.62
10:AJ:6:ILE:O	10:AJ:71:LEU:HD12	2.00	0.62
17:AQ:94:ASN:O	17:AQ:97:SER:OG	2.16	0.62
19:AS:11:VAL:CG2	19:AS:16:LEU:HD11	2.30	0.62
35:BA:1366:A:H2'	35:BA:1367:A:C8	2.34	0.62
35:BA:1754:C:H5'	50:BT:101:PHE:CE1	2.34	0.62
35:BA:1925:C:C2'	35:BA:1926:U:H5'	2.30	0.62
35:BA:2056:G:N2	35:BA:2057:A:H1'	2.15	0.62
35:BA:2795:G:C2	35:BA:2799:C:H5'	2.35	0.62
35:BA:851:U:H2'	35:BA:852:G:C8	2.31	0.62
35:BA:853:G:H2'	35:BA:854:G:H8	1.64	0.62
36:BB:20:C:C2'	36:BB:21:G:H5''	2.27	0.62
38:BD:186:HIS:HB3	38:BD:189:CYS:SG	2.39	0.62
38:BD:260:ARG:HH11	38:BD:260:ARG:HG2	1.64	0.62
41:BG:168:GLU:O	41:BG:171:ALA:HB3	1.99	0.62
42:BH:146:ALA:O	42:BH:147:ASN:C	2.38	0.62
44:BN:16:ILE:HG23	44:BN:54:VAL:CG2	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:111:PHE:HB3	45:BO:114:ILE:CD1	2.27	0.62
45:BO:26:LYS:HB2	45:BO:30:ALA:HB2	1.81	0.62
39:BE:11:MET:H	50:BT:8:LYS:HZ1	1.46	0.62
52:BV:14:VAL:HG12	52:BV:15:GLU:N	2.14	0.62
54:BX:32:PRO:HD3	54:BX:72:LYS:NZ	2.15	0.62
1:CA:16:A:N1	1:CA:919:A:H2	1.98	0.62
1:CA:375:U:H2'	1:CA:376:G:C8	2.34	0.62
1:CA:892:A:H2'	1:CA:893:C:H6	1.65	0.62
2:CB:97:TRP:HH2	2:CB:176:GLU:HB2	1.64	0.62
5:CE:91:LEU:HG	5:CE:120:THR:HG22	1.82	0.62
5:CE:12:LEU:HD13	5:CE:12:LEU:H	1.65	0.62
9:CI:116:LYS:O	9:CI:118:LYS:N	2.33	0.62
9:CI:4:TYR:CD1	9:CI:4:TYR:N	2.68	0.62
11:CK:62:GLN:O	11:CK:64:ALA:N	2.32	0.62
22:CV:28:G:O2'	22:CV:29:G:H5'	2.00	0.62
34:D8:32:LEU:O	34:D8:34:TRP:N	2.32	0.62
34:D8:34:TRP:O	34:D8:35:GLN:HB2	1.98	0.62
31:D5:4:HIS:O	35:DA:2056:G:N2	2.33	0.62
35:DA:2538:C:C2'	35:DA:2539:C:H5'	2.29	0.62
35:DA:2723:C:H2'	35:DA:2724:C:H5'	1.82	0.62
35:DA:2887:U:O2'	35:DA:2888:C:H5'	1.99	0.62
35:DA:2893:G:C5'	35:DA:2894:G:H5'	2.13	0.62
35:DA:744:G:OP1	39:DE:132:HIS:HB3	2.00	0.62
35:DA:76:C:H2'	35:DA:77:C:H6	1.64	0.62
45:DO:68:GLU:HB3	45:DO:78:ARG:HB2	1.82	0.62
47:DQ:35:VAL:HG12	47:DQ:130:LYS:HB3	1.80	0.62
50:DT:28:VAL:HG21	50:DT:46:GLU:HA	1.81	0.62
1:AA:59:A:H3'	1:AA:331:G:H22	1.64	0.61
1:AA:715:A:H2'	1:AA:716:A:H8	1.64	0.61
1:AA:893:C:H2'	1:AA:894:G:C8	2.35	0.61
1:AA:946:A:H2'	1:AA:947:G:H8	1.64	0.61
3:AC:107:GLN:O	3:AC:108:ASN:HB2	2.00	0.61
4:AD:23:GLY:HA3	4:AD:112:VAL:HG22	1.82	0.61
7:AG:150:ALA:C	7:AG:152:ALA:H	2.02	0.61
12:AL:32:PHE:HB3	12:AL:84:LEU:CD2	2.29	0.61
18:AR:81:PHE:O	18:AR:82:THR:HB	1.98	0.61
20:AT:14:LYS:HA	20:AT:17:ARG:HH21	1.64	0.61
23:AW:7:G:H5'	23:AW:8:U:C5	2.35	0.61
25:AY:103:ILE:O	25:AY:103:ILE:HG13	2.00	0.61
27:B1:41:ARG:HG3	27:B1:41:ARG:HH11	1.63	0.61
35:BA:990:A:N6	35:BA:1186:G:H1'	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1192:G:O2'	35:BA:1193:G:H5'	1.99	0.61
35:BA:2673:G:O2'	35:BA:2674:G:H5'	1.99	0.61
35:BA:2875:C:H4'	50:BT:5:ALA:HB2	1.81	0.61
35:BA:491:G:O2'	35:BA:492:A:H5'	2.00	0.61
35:BA:962:G:O2'	35:BA:963:U:H5'	2.00	0.61
37:BC:184:LYS:C	37:BC:186:ALA:H	2.03	0.61
38:BD:145:VAL:HG12	38:BD:146:GLU:N	2.15	0.61
38:BD:224:ALA:HB2	38:BD:233:HIS:HB3	1.80	0.61
41:BG:69:ALA:CB	41:BG:91:ARG:HE	2.13	0.61
43:BI:88:ILE:CG2	43:BI:89:TYR:H	2.12	0.61
45:BO:80:ASP:O	45:BO:81:ASP:HB3	1.99	0.61
48:BR:28:LEU:HA	48:BR:34:ILE:CG1	2.30	0.61
52:BV:96:ILE:CG2	52:BV:97:LYS:N	2.62	0.61
54:BX:60:ARG:HG2	54:BX:74:PRO:CD	2.24	0.61
55:BY:27:VAL:HG12	55:BY:29:GLU:OE1	1.99	0.61
1:CA:1355:G:H2'	1:CA:1356:G:H8	1.63	0.61
2:CB:54:THR:O	2:CB:58:ILE:HG13	2.00	0.61
4:CD:61:LYS:HE3	4:CD:207:TYR:OH	2.00	0.61
5:CE:87:SER:OG	5:CE:125:SER:HB3	2.00	0.61
12:CL:6:THR:HG22	12:CL:9:GLN:HE21	1.65	0.61
27:D1:47:GLN:HB3	27:D1:64:ALA:CB	2.30	0.61
32:D6:42:TRP:HA	32:D6:42:TRP:CE3	2.34	0.61
35:DA:1344:G:H4'	35:DA:1384:A:N7	2.15	0.61
35:DA:1484:G:H3'	35:DA:1485:G:H5''	1.81	0.61
35:DA:2830:G:H5'	39:DE:58:ARG:NH2	2.14	0.61
35:DA:285:C:H3'	35:DA:286:C:H5''	1.82	0.61
35:DA:332:A:H4'	35:DA:333:G:OP1	1.99	0.61
27:D1:47:GLN:N	35:DA:397:G:OP1	2.23	0.61
35:DA:93:G:H2'	35:DA:94:C:H6	1.64	0.61
35:DA:94(A):G:H2'	35:DA:95:G:C5'	2.26	0.61
35:DA:970:C:H2'	35:DA:971:C:H6	1.65	0.61
40:DF:178:PRO:HG2	40:DF:179:GLU:N	2.14	0.61
44:DN:58:ASP:O	44:DN:60:ILE:N	2.30	0.61
48:DR:103:ARG:NH1	48:DR:103:ARG:HG2	2.15	0.61
50:DT:91:ARG:O	50:DT:93:ARG:N	2.32	0.61
51:DU:106:PHE:O	51:DU:110:VAL:HG23	2.00	0.61
53:DW:37:ARG:HG3	53:DW:37:ARG:HH11	1.65	0.61
1:AA:1162:C:H2'	1:AA:1163:C:C6	2.35	0.61
1:AA:1242:C:H2'	1:AA:1243:C:C6	2.34	0.61
1:AA:1523:G:C6	1:AA:1524:C:C4	2.88	0.61
1:AA:393:A:OP2	16:AP:12:LYS:HD3	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:402:G:H1'	1:AA:620:C:H42	1.65	0.61
1:AA:524:G:H2'	1:AA:525:C:C6	2.36	0.61
1:AA:628:G:O2'	1:AA:629:G:H5'	2.00	0.61
10:AJ:32:ALA:HB1	10:AJ:75:ILE:CD1	2.29	0.61
10:AJ:39:PRO:HA	10:AJ:70:ARG:HA	1.81	0.61
25:AY:63:PRO:HB2	25:AY:64:ARG:CZ	2.29	0.61
25:AY:9:GLU:O	25:AY:13:HIS:HB2	1.98	0.61
33:B7:17:GLY:O	33:B7:18:PHE:C	2.38	0.61
34:B8:25:MET:HG3	46:BP:62:LEU:HD21	1.81	0.61
35:BA:1388:G:H1	35:BA:1399:C:H42	1.48	0.61
35:BA:1387:C:H5'	35:BA:1469:A:H4'	1.82	0.61
35:BA:1830:C:H42	35:BA:1975:G:H1	1.48	0.61
35:BA:2707:G:H2'	35:BA:2708:G:C8	2.35	0.61
35:BA:2682:U:O4	35:BA:2728:U:H1'	2.01	0.61
38:BD:109:ASP:HB3	38:BD:195:ALA:HB3	1.81	0.61
39:BE:116:VAL:HG21	39:BE:122:PHE:CG	2.35	0.61
40:BF:160:ASN:HB3	40:BF:163:VAL:HG23	1.82	0.61
40:BF:202:PHE:C	40:BF:202:PHE:HD1	2.03	0.61
40:BF:28:ILE:HD13	40:BF:28:ILE:N	2.14	0.61
43:BI:87:LYS:NZ	43:BI:121:LYS:HG2	2.14	0.61
44:BN:15:LEU:HB3	44:BN:136:GLU:HA	1.82	0.61
49:BS:17:ARG:HG3	49:BS:18:ILE:HD12	1.80	0.61
55:BY:76:CYS:SG	55:BY:77:PRO:CD	2.88	0.61
1:CA:1075:C:OP1	2:CB:179:LYS:HE2	2.00	0.61
1:CA:1313:U:OP1	19:CS:6:LYS:HG3	2.00	0.61
1:CA:1410:G:H2'	1:CA:1411:C:C6	2.36	0.61
1:CA:1510:U:H2'	1:CA:1511:G:C8	2.36	0.61
1:CA:579:G:H5'	1:CA:728:A:H1'	1.82	0.61
3:CC:141:VAL:O	3:CC:141:VAL:HG12	2.00	0.61
3:CC:70:VAL:HG12	3:CC:72:LYS:H	1.64	0.61
12:CL:22:SER:O	12:CL:24:VAL:N	2.33	0.61
13:CM:19:LEU:CA	13:CM:22:ILE:HD13	2.30	0.61
17:CQ:69:LYS:C	17:CQ:70:ARG:HD2	2.20	0.61
19:CS:11:VAL:CG2	19:CS:16:LEU:HD11	2.30	0.61
26:D0:25:ARG:HG3	26:D0:29:GLN:NE2	2.14	0.61
35:DA:1440:G:H2'	35:DA:1441:G:H8	1.64	0.61
35:DA:1794:U:O4'	35:DA:1900:A:C2	2.53	0.61
35:DA:2078:C:H2'	35:DA:2079:U:C6	2.34	0.61
35:DA:2225:A:H4'	35:DA:2226:C:H5'	1.83	0.61
35:DA:2308:G:H8	35:DA:2309:A:H3'	1.65	0.61
35:DA:2389:G:H5''	35:DA:2390:U:O4'	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2684:U:P	50:DT:53:ARG:HH21	2.23	0.61
35:DA:2725:A:O2'	35:DA:2726:U:H2'	2.00	0.61
35:DA:2795:G:N2	35:DA:2799:C:H5'	2.14	0.61
35:DA:2820:A:O3'	48:DR:2:ARG:NH2	2.33	0.61
35:DA:61:G:H1	35:DA:94:C:H42	1.48	0.61
41:DG:22:ARG:O	41:DG:22:ARG:HD3	1.99	0.61
41:DG:46:ALA:O	41:DG:51:ARG:HG3	1.99	0.61
41:DG:86:MET:N	41:DG:87:PRO:HD3	2.15	0.61
42:DH:44:VAL:HG12	42:DH:45:VAL:N	2.14	0.61
44:DN:10:GLU:CG	44:DN:11:PRO:HD2	2.30	0.61
44:DN:17:ASP:C	44:DN:19:GLU:N	2.53	0.61
44:DN:58:ASP:O	44:DN:60:ILE:HG13	2.00	0.61
44:DN:62:VAL:HG22	44:DN:66:LYS:CG	2.30	0.61
45:DO:105:GLU:HA	45:DO:108:GLU:CG	2.30	0.61
48:DR:13:HIS:CE1	48:DR:16:HIS:HB2	2.35	0.61
48:DR:84:ALA:HB3	48:DR:85:PRO:HD3	1.80	0.61
50:DT:48:ILE:O	50:DT:63:VAL:HG12	2.00	0.61
51:DU:33:ARG:O	51:DU:34:LYS:C	2.38	0.61
28:D2:26:ARG:NH2	54:DX:6:ASP:HA	2.13	0.61
55:DY:95:LYS:HG2	55:DY:100:ALA:CA	2.20	0.61
56:DZ:31:ARG:CZ	56:DZ:94:GLU:OE2	2.48	0.61
1:AA:155:C:H2'	1:AA:156:G:C8	2.36	0.61
1:AA:184:G:C4'	1:AA:224:C:H4'	2.31	0.61
2:AB:97:TRP:HH2	2:AB:176:GLU:HB2	1.65	0.61
6:AF:76:ALA:HB1	6:AF:80:ARG:HH21	1.64	0.61
7:AG:105:VAL:HG12	7:AG:109:ASN:ND2	2.15	0.61
8:AH:12:ARG:HH11	8:AH:26:VAL:HA	1.65	0.61
8:AH:18:ARG:N	8:AH:78:GLN:HE22	1.98	0.61
16:AP:48:TRP:HE3	16:AP:49:LEU:H	1.48	0.61
35:BA:1190:G:H4'	46:BP:35:HIS:CB	2.29	0.61
35:BA:1547:C:O2'	35:BA:1548:C:H5'	2.00	0.61
27:B1:37:ILE:HG21	35:BA:2079:U:O3'	1.99	0.61
26:B0:14:ARG:HD2	35:BA:2279:G:O6	2.00	0.61
35:BA:2364:C:H2'	35:BA:2365:G:O4'	1.99	0.61
35:BA:2723:C:O2'	35:BA:2724:C:H5'	2.00	0.61
35:BA:622:G:O2'	35:BA:623:G:H5'	2.01	0.61
37:BC:211:SER:HA	37:BC:220:PRO:HA	1.82	0.61
39:BE:116:VAL:O	39:BE:117:MET:CB	2.47	0.61
39:BE:5:LEU:HD22	39:BE:197:ILE:HG22	1.81	0.61
40:BF:67:GLN:CG	40:BF:67:GLN:O	2.41	0.61
41:BG:51:ARG:HA	41:BG:51:ARG:HE	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BH:87:LEU:N	42:BH:131:VAL:O	2.34	0.61
43:BI:95:LYS:O	43:BI:99:GLU:HB2	2.00	0.61
52:BV:39:LEU:HD21	52:BV:53:GLU:HA	1.82	0.61
35:BA:1614:A:N6	53:BW:93:ALA:CB	2.63	0.61
54:BX:65:ARG:CZ	54:BX:66:LEU:N	2.61	0.61
56:BZ:142:SER:H	56:BZ:144:LEU:HD23	1.64	0.61
56:BZ:17:ALA:HA	56:BZ:20:ARG:CD	2.29	0.61
1:CA:1010:G:H22	1:CA:1020:U:H1'	1.64	0.61
1:CA:1160:G:OP1	2:CB:132:LYS:HE3	2.01	0.61
1:CA:1405:G:O2'	1:CA:1406:U:H5'	1.99	0.61
1:CA:1517:G:C8	35:DA:1920:C:OP1	2.53	0.61
1:CA:519:C:H2'	1:CA:520:A:C8	2.36	0.61
4:CD:60:GLU:HG2	4:CD:202:LEU:HB2	1.82	0.61
5:CE:15:ARG:HG3	5:CE:28:PHE:CE2	2.35	0.61
5:CE:57:LYS:HE2	5:CE:61:TYR:HE2	1.66	0.61
12:CL:60:LEU:HD23	12:CL:64:TYR:O	2.00	0.61
16:CP:49:LEU:HD12	16:CP:50:LYS:H	1.62	0.61
34:D8:26:LYS:NZ	34:D8:47:LYS:HD3	2.15	0.61
35:DA:1210:A:H4'	35:DA:1211:U:O5'	2.00	0.61
35:DA:1771:C:H2'	35:DA:1772:G:H8	1.65	0.61
35:DA:2052:G:N3	39:DE:149:ARG:HA	2.16	0.61
35:DA:2165:G:N3	35:DA:2165:G:H2'	2.15	0.61
36:DB:83:G:O2'	36:DB:84:C:H5'	2.01	0.61
38:DD:177:LEU:HD12	38:DD:181:GLU:CG	2.30	0.61
38:DD:68:LYS:HB2	38:DD:70:TRP:CH2	2.34	0.61
39:DE:93:VAL:C	39:DE:95:ILE:H	2.03	0.61
41:DG:38:VAL:HB	41:DG:158:ALA:HB3	1.83	0.61
43:DI:14:ASP:HB2	43:DI:17:GLN:OE1	2.01	0.61
43:DI:4:ILE:HA	43:DI:17:GLN:O	2.00	0.61
45:DO:71:ARG:HH11	45:DO:71:ARG:HG3	1.65	0.61
45:DO:93:PRO:C	45:DO:95:GLY:H	2.03	0.61
46:DP:100:LEU:H	46:DP:100:LEU:HD22	1.66	0.61
49:DS:72:ALA:O	49:DS:76:LYS:HG2	2.00	0.61
39:DE:11:MET:H	50:DT:8:LYS:HZ1	1.47	0.61
50:DT:50:ILE:CG2	50:DT:99:LEU:HD12	2.23	0.61
52:DV:61:VAL:CG2	52:DV:100:ARG:HG2	2.31	0.61
53:DW:69:LEU:H	53:DW:69:LEU:HD12	1.64	0.61
1:AA:1003:G:H2'	1:AA:1004:A:O4'	1.99	0.61
1:AA:1355:G:H2'	1:AA:1356:G:H8	1.65	0.61
1:AA:1478:C:C2	1:AA:1479:C:C5	2.88	0.61
3:AC:164:ARG:CB	3:AC:164:ARG:HH11	2.06	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:40:ALA:HA	7:AG:43:PHE:HB3	1.81	0.61
5:AE:78:HIS:CD2	8:AH:104:ARG:HE	2.19	0.61
9:AI:4:TYR:CD1	9:AI:4:TYR:N	2.68	0.61
11:AK:23:ALA:HB1	11:AK:91:ARG:HG2	1.81	0.61
13:AM:78:ILE:HA	13:AM:81:LEU:CD1	2.30	0.61
20:AT:57:ARG:HH11	20:AT:57:ARG:HB2	1.65	0.61
35:BA:1192:G:C2'	35:BA:1193:G:H5'	2.30	0.61
35:BA:2037:G:H2'	35:BA:2038:G:C8	2.35	0.61
35:BA:648:G:C4'	35:BA:2351:G:H5''	2.31	0.61
35:BA:2609:U:OP1	35:BA:2609:U:H4'	2.01	0.61
35:BA:2787:C:C2	39:BE:61:ARG:HD3	2.36	0.61
36:BB:31:C:H2'	36:BB:53:A:H61	1.66	0.61
37:BC:213:TYR:CB	37:BC:218:MET:HA	2.30	0.61
38:BD:181:GLU:HA	38:BD:272:ALA:O	2.00	0.61
42:BH:136:ILE:O	42:BH:137:ASP:HB2	1.99	0.61
46:BP:99:LEU:HD12	46:BP:102:ARG:HD2	1.82	0.61
46:BP:38:GLN:HG3	46:BP:39:LYS:N	2.06	0.61
47:BQ:53:ALA:HA	47:BQ:56:ARG:HB3	1.83	0.61
51:BU:40:PHE:N	51:BU:40:PHE:CD2	2.68	0.61
51:BU:55:ARG:HA	51:BU:58:ARG:HD2	1.82	0.61
51:BU:66:ASN:OD1	51:BU:76:TYR:N	2.33	0.61
54:BX:36:LYS:HZ1	54:BX:39:ILE:HA	1.65	0.61
55:BY:37:VAL:HG22	55:BY:67:LEU:O	2.00	0.61
1:CA:1088:G:H2'	1:CA:1089:G:C8	2.33	0.61
1:CA:1408:A:O2'	35:DA:1916:A:N6	2.33	0.61
1:CA:471:G:H2'	1:CA:472:A:C8	2.34	0.61
1:CA:473:G:H2'	1:CA:474:G:H8	1.66	0.61
2:CB:72:GLY:HA3	2:CB:165:VAL:HG22	1.82	0.61
2:CB:212:GLN:CG	2:CB:235:SER:HB2	2.30	0.61
5:CE:152:ARG:HB3	8:CH:43:GLY:HA3	1.82	0.61
6:CF:61:LEU:O	6:CF:62:TRP:HB2	1.99	0.61
8:CH:103:VAL:HG12	8:CH:108:GLY:HA3	1.83	0.61
9:CI:26:VAL:HG13	9:CI:63:ILE:CD1	2.30	0.61
12:CL:89:ARG:HH11	12:CL:90:VAL:N	1.98	0.61
12:CL:90:VAL:HG11	12:CL:93:LEU:HG	1.81	0.61
17:CQ:94:ASN:O	17:CQ:97:SER:OG	2.19	0.61
19:CS:6:LYS:CD	19:CS:7:LYS:HD2	2.30	0.61
24:CX:19:U:H2'	24:CX:19:U:O2	1.99	0.61
35:DA:1388:G:H1	35:DA:1399:C:H42	1.47	0.61
35:DA:1509(A):A:H2'	35:DA:1509(B):A:C8	2.35	0.61
35:DA:2033:A:H4'	35:DA:2034:U:OP1	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2498:C:HO2'	35:DA:2499:C:H5'	1.66	0.61
35:DA:74:A:H4'	35:DA:75:G:O5'	1.99	0.61
35:DA:869:G:H2'	35:DA:870:A:H8	1.65	0.61
35:DA:1803:A:O2'	38:DD:259:THR:HG21	1.99	0.61
38:DD:27:THR:CG2	38:DD:28:GLU:H	1.91	0.61
39:DE:134:ILE:HD13	39:DE:134:ILE:N	2.15	0.61
39:DE:48:GLN:HG2	39:DE:78:LEU:HD12	1.82	0.61
41:DG:114:ILE:CB	41:DG:117:PHE:HB2	2.30	0.61
41:DG:135:LEU:CD2	41:DG:140:ILE:HD11	2.28	0.61
41:DG:96:ARG:HG3	41:DG:97:ASP:H	1.64	0.61
43:DI:51:ILE:HG22	43:DI:52:ARG:N	2.14	0.61
44:DN:120:LEU:HD13	44:DN:120:LEU:C	2.21	0.61
45:DO:102:VAL:HB	45:DO:106:LEU:HD12	1.81	0.61
54:DX:77:LYS:CE	54:DX:78:LYS:HG3	2.30	0.61
54:DX:89:ILE:HA	54:DX:92:LEU:HD12	1.82	0.61
1:AA:1385:G:O2'	1:AA:1386:G:H5'	2.00	0.61
1:AA:253:U:H2'	1:AA:254:G:C8	2.31	0.61
1:AA:383:A:H2'	1:AA:384:G:C5'	2.29	0.61
1:AA:490:G:H2'	1:AA:491:G:C8	2.34	0.61
1:AA:571:U:H5''	1:AA:819:A:C2	2.35	0.61
1:AA:786:G:H1	1:AA:796:C:N4	1.98	0.61
2:AB:33:TYR:CD1	2:AB:43:ASP:HA	2.34	0.61
6:AF:91:VAL:HG11	18:AR:72:ARG:NH1	2.15	0.61
12:AL:22:SER:O	12:AL:24:VAL:N	2.33	0.61
12:AL:47:LYS:CB	12:AL:48:PRO:CD	2.79	0.61
18:AR:37:VAL:CG2	18:AR:38:GLU:H	2.06	0.61
19:AS:40:ILE:HD13	19:AS:62:ILE:HD13	1.83	0.61
23:AW:60:A:C2'	23:AW:61:U:H5'	2.30	0.61
25:AY:52:LEU:C	25:AY:54:GLN:H	2.03	0.61
25:AY:80:GLU:O	25:AY:82:ALA:N	2.33	0.61
28:B2:22:GLU:HA	28:B2:25:VAL:CG1	2.31	0.61
29:B3:36:VAL:O	29:B3:37:LEU:HD23	2.00	0.61
35:BA:1153:C:H2'	35:BA:1154:G:O4'	2.00	0.61
35:BA:143:G:H2'	35:BA:143(A):C:C6	2.35	0.61
35:BA:1660:C:H5'	35:BA:2712(A):A:H61	1.63	0.61
35:BA:2111:C:H5'	35:BA:2112:G:OP1	2.00	0.61
35:BA:2078:C:H1'	35:BA:2434:A:N3	2.16	0.61
35:BA:285:C:H3'	35:BA:286:C:H5''	1.82	0.61
35:BA:506:G:H4'	35:BA:509:C:O2	2.01	0.61
35:BA:595:C:H2'	35:BA:596:G:C8	2.36	0.61
35:BA:260:G:H1'	35:BA:621:A:H1'	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:796:C:H2'	35:BA:797:C:C6	2.36	0.61
38:BD:10:THR:HG23	38:BD:13:ARG:HB3	1.82	0.61
38:BD:166:GLN:N	38:BD:166:GLN:HE21	1.98	0.61
38:BD:231:HIS:CD2	38:BD:249:PRO:HG3	2.36	0.61
39:BE:103:ASP:OD2	39:BE:201:THR:HA	2.00	0.61
39:BE:197:ILE:CD1	39:BE:199:ARG:HH22	2.14	0.61
40:BF:139:PHE:CB	40:BF:166:ALA:HB1	2.31	0.61
40:BF:202:PHE:CD1	40:BF:202:PHE:C	2.73	0.61
41:BG:77:ILE:HD13	41:BG:80:PHE:O	2.01	0.61
42:BH:86:GLU:HA	42:BH:132:ARG:HB3	1.81	0.61
43:BI:91:SER:H	43:BI:121:LYS:CE	2.13	0.61
44:BN:78:TYR:H	44:BN:79:PRO:HD2	1.65	0.61
45:BO:71:ARG:HH11	45:BO:71:ARG:HG3	1.64	0.61
45:BO:76:ALA:HB3	50:BT:75:ILE:CB	2.30	0.61
35:BA:598:G:H5'	46:BP:15:ARG:CD	2.31	0.61
49:BS:38:GLN:C	49:BS:39:ILE:HD12	2.20	0.61
51:BU:55:ARG:HA	51:BU:58:ARG:CD	2.30	0.61
54:BX:82:GLN:CG	54:BX:83:VAL:N	2.61	0.61
1:CA:17:U:O2'	1:CA:18:C:H5'	2.00	0.61
1:CA:585:G:H4'	12:CL:8:ASN:ND2	2.10	0.61
1:CA:640:A:H4'	8:CH:116:LYS:NZ	2.15	0.61
1:CA:673:G:H5''	6:CF:87:ARG:HH11	1.63	0.61
2:CB:36:ARG:NH1	2:CB:37:ASN:HB2	2.16	0.61
8:CH:119:LEU:HD23	8:CH:119:LEU:H	1.63	0.61
8:CH:64:LYS:O	8:CH:79:VAL:HG21	2.01	0.61
22:CV:40:C:O2'	22:CV:41:C:H5'	2.01	0.61
27:D1:10:LYS:C	27:D1:11:ARG:O	2.36	0.61
35:DA:1665:A:H2'	35:DA:1666:G:O4'	2.00	0.61
35:DA:1814:G:H2'	35:DA:1815:A:N7	2.15	0.61
35:DA:2203:U:C4'	38:DD:151:LYS:HE3	2.30	0.61
35:DA:581:C:O2'	35:DA:582:G:H5'	2.00	0.61
38:DD:145:VAL:HB	38:DD:155:LEU:HB2	1.83	0.61
35:DA:2575:C:H5'	39:DE:144:ARG:HG2	1.82	0.61
39:DE:92:THR:HB	39:DE:94:GLU:OE1	2.00	0.61
40:DF:167:ALA:HB1	40:DF:173:VAL:HG11	1.82	0.61
40:DF:178:PRO:CG	40:DF:179:GLU:H	2.11	0.61
35:DA:675:A:H4'	40:DF:67:GLN:OE1	2.00	0.61
41:DG:120:LEU:O	41:DG:180:PHE:HA	2.01	0.61
47:DQ:29:PHE:HB3	56:DZ:118:GLN:OE1	2.00	0.61
50:DT:91:ARG:O	50:DT:117:ASP:HB2	2.01	0.61
55:DY:38:ILE:HG22	55:DY:39:VAL:N	2.14	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:86:VAL:HG12	56:DZ:87:ASP:N	2.14	0.61
1:AA:17:U:O4'	1:AA:1080:A:H1'	2.00	0.61
1:AA:135:C:H2'	1:AA:136:C:H5'	1.82	0.61
1:AA:1463:C:H2'	1:AA:1464:G:O4'	2.01	0.61
1:AA:17:U:O2'	1:AA:18:C:H5'	2.00	0.61
1:AA:818:G:C3'	1:AA:819:A:H5''	2.31	0.61
2:AB:103:THR:HA	2:AB:180:LEU:HD11	1.82	0.61
2:AB:80:ILE:HG21	2:AB:208:ILE:HG23	1.83	0.61
3:AC:119:ARG:NH2	3:AC:140:ARG:CZ	2.63	0.61
13:AM:115:LYS:O	13:AM:117:VAL:HG23	2.00	0.61
15:AO:53:HIS:HE1	15:AO:57:LEU:HD21	1.66	0.61
17:AQ:67:LYS:HA	17:AQ:70:ARG:HH12	1.64	0.61
23:AW:45:A:H2'	23:AW:46:G:C8	2.36	0.61
35:BA:1523:U:H2'	35:BA:1524:G:C8	2.36	0.61
35:BA:2179:C:H1'	37:BC:171:ILE:CB	2.30	0.61
35:BA:2683:C:H5''	50:BT:53:ARG:NH2	2.14	0.61
35:BA:696:G:O2'	35:BA:697:C:H5'	1.99	0.61
38:BD:222:ARG:O	38:BD:223:GLY:C	2.37	0.61
44:BN:1:MET:C	44:BN:2:LYS:HD2	2.21	0.61
45:BO:102:VAL:HB	45:BO:106:LEU:HD12	1.83	0.61
46:BP:46:LYS:HB3	46:BP:52:GLU:HG2	1.81	0.61
47:BQ:60:ARG:C	47:BQ:60:ARG:HD3	2.20	0.61
49:BS:88:ASP:OD2	49:BS:89:ARG:N	2.29	0.61
52:BV:2:PHE:HB3	52:BV:42:GLY:CA	2.26	0.61
56:BZ:57:ILE:HG22	56:BZ:59:LEU:HG	1.81	0.61
1:CA:1152:A:H5''	10:CJ:13:HIS:HD2	1.66	0.61
1:CA:1465:C:H2'	1:CA:1466:C:C6	2.36	0.61
1:CA:711:G:O2'	1:CA:712:A:H5'	2.01	0.61
3:CC:164:ARG:CB	3:CC:164:ARG:HH11	2.07	0.61
8:CH:26:VAL:O	8:CH:59:LEU:N	2.33	0.61
27:D1:42:GLN:HG2	27:D1:43:TYR:N	2.15	0.61
28:D2:15:LYS:O	28:D2:18:PRO:HD2	2.00	0.61
33:D7:29:LYS:O	33:D7:32:LYS:HB3	2.01	0.61
34:D8:25:MET:HB2	46:DP:62:LEU:HD11	1.82	0.61
35:DA:1366:A:H2'	35:DA:1367:A:H8	1.65	0.61
35:DA:1854:A:H5'	35:DA:1855:G:OP2	1.99	0.61
35:DA:1925:C:C2'	35:DA:1926:U:H5'	2.30	0.61
35:DA:491:G:O2'	35:DA:492:A:H5'	1.99	0.61
35:DA:499:U:O2'	35:DA:500:G:H5'	2.00	0.61
35:DA:595:C:H2'	35:DA:596:G:C8	2.36	0.61
42:DH:88:LEU:O	42:DH:89:ILE:HG23	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DI:64:GLU:OE1	43:DI:67:ARG:HB2	2.01	0.61
44:DN:32:THR:CG2	44:DN:37:LYS:HB3	2.31	0.61
46:DP:130:PHE:N	46:DP:130:PHE:HD2	1.98	0.61
47:DQ:43:THR:OG1	47:DQ:45:GLN:HB2	2.01	0.61
47:DQ:76:LYS:H	47:DQ:88:GLY:HA2	1.64	0.61
51:DU:31:SER:C	51:DU:33:ARG:H	2.03	0.61
56:DZ:146:ILE:HA	56:DZ:174:VAL:CB	2.29	0.61
56:DZ:24:LEU:C	56:DZ:24:LEU:HD12	2.21	0.61
56:DZ:97:GLU:HA	56:DZ:126:VAL:O	2.00	0.61
1:AA:1080:A:H5'	5:AE:14:ARG:NH2	2.14	0.61
1:AA:620:C:H2'	1:AA:621:A:O4'	2.01	0.61
1:AA:685:G:N2	1:AA:686:U:H3	1.99	0.61
2:AB:178:ARG:HH22	2:AB:196:LEU:HA	1.65	0.61
2:AB:185:ILE:HG22	2:AB:199:TYR:CD1	2.35	0.61
6:AF:37:VAL:HG13	6:AF:65:VAL:CG1	2.31	0.61
6:AF:82:ARG:HB3	6:AF:82:ARG:NH1	2.15	0.61
7:AG:58:PRO:HA	7:AG:61:VAL:CG2	2.31	0.61
8:AH:103:VAL:HG21	8:AH:109:ILE:C	2.21	0.61
9:AI:3:GLN:HG2	9:AI:20:ARG:HH21	1.65	0.61
12:AL:119:LYS:HB2	12:AL:120:TYR:HD1	1.64	0.61
18:AR:43:PHE:C	18:AR:44:LEU:HD12	2.21	0.61
6:AF:89:MET:SD	18:AR:76:LEU:HD21	2.41	0.61
26:B0:25:ARG:HG3	26:B0:29:GLN:NE2	2.15	0.61
29:B3:41:PRO:HD3	29:B3:44:ARG:NH1	2.16	0.61
35:BA:1821:A:H2'	35:BA:1822:G:C8	2.35	0.61
35:BA:2505:G:H2'	35:BA:2576:G:O6	2.01	0.61
35:BA:2830:G:H5'	39:BE:58:ARG:NH2	2.16	0.61
35:BA:426:C:O2'	35:BA:427:U:H5'	2.01	0.61
39:BE:141:ILE:HD13	39:BE:141:ILE:N	2.16	0.61
35:BA:2619:C:H5''	39:BE:152:LYS:HA	1.83	0.61
45:BO:11:ALA:HB1	45:BO:99:PHE:O	2.01	0.61
45:BO:1:MET:HG3	45:BO:32:TYR:HD2	1.59	0.61
46:BP:85:LEU:CD2	46:BP:85:LEU:H	2.09	0.61
35:BA:1286:A:OP1	48:BR:105:ARG:HD2	2.01	0.61
48:BR:9:LYS:HE3	48:BR:43:GLU:OE2	2.01	0.61
50:BT:48:ILE:O	50:BT:63:VAL:HG12	2.01	0.61
51:BU:105:VAL:O	51:BU:109:LEU:HG	2.01	0.61
51:BU:90:VAL:HG22	52:BV:39:LEU:CD1	2.30	0.61
53:BW:5:ALA:HB2	53:BW:54:ALA:HB2	1.83	0.61
54:BX:89:ILE:O	54:BX:89:ILE:HG22	1.99	0.61
1:CA:431:A:H2'	1:CA:432:A:C8	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:33:TYR:HB2	2:CB:41:ILE:HG22	1.81	0.61
5:CE:139:LEU:HA	5:CE:142:LEU:CD1	2.30	0.61
9:CI:7:THR:HB	9:CI:83:ARG:NH1	2.14	0.61
13:CM:117:VAL:HG12	13:CM:118:ALA:N	2.16	0.61
18:CR:38:GLU:O	18:CR:41:LYS:HB3	1.99	0.61
20:CT:31:SER:O	20:CT:34:LYS:HB2	2.00	0.61
20:CT:64:ASP:C	20:CT:66:ALA:H	2.02	0.61
35:DA:1215:G:H2'	35:DA:1216:G:H8	1.66	0.61
35:DA:2260:C:O2'	35:DA:2261:C:H5'	1.99	0.61
35:DA:2707:G:H5''	48:DR:68:ARG:HH21	1.66	0.61
36:DB:31:C:H2'	36:DB:53:A:H61	1.66	0.61
41:DG:111:LEU:CA	41:DG:114:ILE:HD11	2.29	0.61
43:DI:77:LEU:CB	43:DI:140:LEU:HD13	2.22	0.61
45:DO:42:SER:HA	45:DO:56:ASP:O	1.99	0.61
47:DQ:34:LEU:CD1	47:DQ:129:THR:HB	2.29	0.61
49:DS:31:SER:OG	49:DS:32:LEU:N	2.32	0.61
49:DS:41:ASP:O	49:DS:45:GLY:HA2	2.00	0.61
49:DS:67:ARG:HE	49:DS:100:ALA:HB3	1.66	0.61
53:DW:26:GLY:O	53:DW:27:LYS:HG3	2.00	0.61
54:DX:39:ILE:O	54:DX:42:ALA:HB3	2.00	0.61
54:DX:30:VAL:HG23	54:DX:76:ARG:HA	1.81	0.61
54:DX:82:GLN:CD	54:DX:83:VAL:H	2.04	0.61
55:DY:19:LYS:HD2	55:DY:20:TYR:CE1	2.36	0.61
56:DZ:98:MET:O	56:DZ:98:MET:HG3	2.00	0.61
1:AA:1056:U:H5'	3:AC:163:ALA:HB2	1.82	0.61
1:AA:291:C:O2'	1:AA:292:G:H5'	2.01	0.61
1:AA:707:C:H4'	11:AK:20:TYR:HD1	1.64	0.61
1:AA:1190:G:H8	3:AC:3:ASN:HD21	1.47	0.61
6:AF:23:LYS:O	6:AF:27:GLN:HG2	2.01	0.61
12:AL:89:ARG:HE	12:AL:91:LYS:HE2	1.66	0.61
13:AM:97:PRO:HB3	13:AM:110:ARG:HD3	1.81	0.61
15:AO:24:SER:O	15:AO:28:GLN:HG3	2.01	0.61
16:AP:72:ARG:C	16:AP:74:LEU:H	2.04	0.61
25:AY:30:THR:C	25:AY:32:ARG:N	2.53	0.61
27:B1:73:LEU:CA	27:B1:76:ARG:HH12	2.14	0.61
32:B6:42:TRP:CE3	32:B6:42:TRP:HA	2.35	0.61
33:B7:13:ALA:O	33:B7:17:GLY:HA3	2.00	0.61
35:BA:1220:A:O2'	35:BA:1221:C:H5''	2.00	0.61
35:BA:1889:A:O2'	35:BA:2087:G:H5'	1.99	0.61
35:BA:1902:C:C5'	38:BD:246:PRO:HD3	2.31	0.61
35:BA:1909:C:H2'	35:BA:1909:C:O2	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2241:A:H2'	35:BA:2242:G:C8	2.35	0.61
35:BA:2308:G:H8	35:BA:2309:A:H3'	1.65	0.61
35:BA:2514:U:H2'	35:BA:2515:C:H6	1.65	0.61
35:BA:542:C:H2'	35:BA:543:C:OP1	2.01	0.61
35:BA:559:G:H22	51:BU:49:HIS:CD2	2.19	0.61
35:BA:648:G:H2'	35:BA:649:G:C8	2.35	0.61
35:BA:792:G:C5'	35:BA:793:A:H5'	2.30	0.61
38:BD:9:TYR:C	38:BD:10:THR:HG22	2.21	0.61
39:BE:38:THR:C	39:BE:40:GLU:H	2.04	0.61
39:BE:64:LYS:C	39:BE:66:HIS:H	2.01	0.61
39:BE:52:LEU:N	39:BE:74:PRO:HB2	2.15	0.61
43:BI:96:ASP:HA	43:BI:99:GLU:HB3	1.82	0.61
46:BP:108:LYS:C	46:BP:110:TYR:H	2.04	0.61
46:BP:56:SER:O	46:BP:57:THR:HB	2.01	0.61
48:BR:103:ARG:HG2	48:BR:103:ARG:NH1	2.13	0.61
50:BT:109:GLU:CA	50:BT:112:ARG:HG3	2.30	0.61
53:BW:69:LEU:HD12	53:BW:69:LEU:H	1.65	0.61
55:BY:13:VAL:HG21	55:BY:28:LYS:HG2	1.82	0.61
1:CA:694:A:H5''	11:CK:53:SER:CB	2.30	0.61
3:CC:153:VAL:HA	3:CC:197:GLY:O	2.01	0.61
3:CC:182:ILE:HG12	3:CC:203:PHE:HA	1.82	0.61
4:CD:23:GLY:HA3	4:CD:112:VAL:HG22	1.82	0.61
5:CE:131:ILE:HD13	5:CE:131:ILE:H	1.64	0.61
7:CG:149:ARG:HD3	11:CK:59:TYR:CE1	2.36	0.61
13:CM:58:GLU:HA	13:CM:58:GLU:OE1	2.01	0.61
17:CQ:5:VAL:HA	17:CQ:59:ILE:O	2.01	0.61
28:D2:49:LYS:O	28:D2:51:ARG:N	2.34	0.61
31:D5:52:TYR:HA	31:D5:56:LYS:NZ	2.15	0.61
35:DA:1014:U:H2'	35:DA:1015:G:H8	1.65	0.61
35:DA:225:A:O2'	35:DA:226:G:H5'	2.01	0.61
35:DA:45:C:H2'	35:DA:47:C:H6	1.66	0.61
38:DD:121:PRO:HB3	38:DD:135:PHE:CE2	2.36	0.61
35:DA:778:G:C5'	38:DD:48:ARG:HD2	2.31	0.61
38:DD:64:ILE:HG23	38:DD:64:ILE:O	2.00	0.61
39:DE:39:PRO:HA	39:DE:43:GLY:CA	2.30	0.61
40:DF:160:ASN:HB3	40:DF:163:VAL:HG23	1.81	0.61
42:DH:40:GLU:HB2	42:DH:41:MET:SD	2.41	0.61
43:DI:10:GLU:O	43:DI:12:LEU:HD23	2.00	0.61
49:DS:85:VAL:HG23	49:DS:106:ARG:HB2	1.82	0.61
54:DX:92:LEU:O	54:DX:93:GLU:HB3	2.01	0.61
56:DZ:58:VAL:HA	56:DZ:67:LEU:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1346:A:H61	1:AA:1374:A:H3'	1.66	0.61
1:AA:414:A:H2'	1:AA:415:A:O4'	2.01	0.61
1:AA:731:G:OP1	1:AA:766:A:H1'	2.00	0.61
2:AB:72:GLY:CA	2:AB:165:VAL:HG22	2.31	0.61
6:AF:18:GLN:O	6:AF:21:LEU:HB2	2.01	0.61
8:AH:85:ARG:HH12	8:AH:134:ILE:HG23	1.66	0.61
25:AY:16:LYS:HA	25:AY:19:GLU:HG3	1.83	0.61
28:B2:41:ILE:O	28:B2:44:LEU:N	2.32	0.61
35:BA:143:G:H2'	35:BA:143(A):C:H6	1.66	0.61
35:BA:2165:G:H2'	35:BA:2165:G:N3	2.16	0.61
35:BA:2348:U:H2'	35:BA:2349:G:C5'	2.23	0.61
35:BA:2460:U:O2'	35:BA:2461:C:H5'	2.00	0.61
35:BA:2677:G:H2'	35:BA:2678:C:C6	2.34	0.61
35:BA:361:G:C2'	35:BA:362:U:H5''	2.31	0.61
35:BA:84:A:N3	35:BA:85:G:H1'	2.15	0.61
36:BB:15:A:H3'	36:BB:16:G:H5'	1.83	0.61
37:BC:82:LYS:HB3	37:BC:86:ALA:HB2	1.82	0.61
35:BA:2512:C:H4'	39:BE:122:PHE:CZ	2.36	0.61
44:BN:43:THR:O	44:BN:46:VAL:N	2.33	0.61
46:BP:92:GLU:HG3	46:BP:93:GLY:H	1.66	0.61
50:BT:28:VAL:O	50:BT:88:ILE:HD11	2.01	0.61
44:BN:42:TRP:HD1	51:BU:63:VAL:HG11	1.66	0.61
52:BV:43:GLU:HB2	52:BV:48:GLY:HA3	1.83	0.61
1:CA:861:G:H2'	1:CA:862:C:H6	1.66	0.61
2:CB:172:ILE:HD12	2:CB:172:ILE:N	2.12	0.61
3:CC:60:ALA:HB3	3:CC:63:ASN:HD21	1.66	0.61
4:CD:108:LEU:CB	4:CD:110:PHE:CE1	2.84	0.61
5:CE:7:GLU:O	5:CE:8:GLU:HB3	1.99	0.61
7:CG:58:PRO:HA	7:CG:61:VAL:CG2	2.31	0.61
8:CH:10:LEU:HD23	8:CH:10:LEU:N	2.16	0.61
11:CK:23:ALA:HB1	11:CK:91:ARG:HG2	1.83	0.61
16:CP:20:VAL:HG23	16:CP:34:GLU:O	2.01	0.61
1:CA:472:A:H1'	16:CP:82:GLN:OE1	2.01	0.61
25:CY:150:SER:HB3	25:CY:153:GLU:HG3	1.81	0.61
28:D2:12:GLU:O	28:D2:14:ARG:HD3	2.01	0.61
35:DA:142:A:H5'	35:DA:142(A):C:OP2	2.01	0.61
35:DA:1496:A:H2'	35:DA:1498:C:C5	2.36	0.61
35:DA:1613:G:H2'	35:DA:1617:C:H42	1.66	0.61
35:DA:1789:A:O2'	35:DA:1790:C:H5'	2.01	0.61
35:DA:202:U:H2'	35:DA:203:C:C6	2.35	0.61
35:DA:2056:G:H2'	35:DA:2056:G:N3	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2128:C:O2'	35:DA:2163:C:OP1	2.17	0.61
35:DA:695:G:N2	35:DA:696:G:H1'	2.15	0.61
35:DA:782:A:H2	38:DD:226:MET:HE2	1.65	0.61
35:DA:820:A:H2'	35:DA:821:A:O4'	2.01	0.61
37:DC:213:TYR:CB	37:DC:218:MET:HA	2.31	0.61
39:DE:199:ARG:HH11	39:DE:199:ARG:HG3	1.66	0.61
40:DF:89:VAL:HG12	40:DF:90:PHE:N	2.16	0.61
41:DG:134:GLY:HA2	41:DG:157:ILE:HG12	1.82	0.61
43:DI:72:LEU:CD1	43:DI:138:ILE:HD11	2.24	0.61
49:DS:23:ARG:HG2	49:DS:24:LEU:H	1.66	0.61
53:DW:110:LYS:HG3	53:DW:111:HIS:ND1	2.16	0.61
1:AA:1078:U:H2'	1:AA:1079:G:C8	2.36	0.61
1:AA:1514:C:H2'	1:AA:1515:C:C6	2.36	0.61
1:AA:710:G:H5''	6:AF:54:LYS:CE	2.31	0.61
4:AD:121:VAL:O	4:AD:134:ASP:HA	2.01	0.61
4:AD:96:LEU:C	4:AD:98:GLU:H	2.02	0.61
6:AF:30:LEU:HB3	6:AF:35:ALA:CB	2.30	0.61
10:AJ:29:ARG:HH11	10:AJ:29:ARG:HG2	1.66	0.61
11:AK:110:ASP:O	18:AR:84:LYS:HD2	2.00	0.61
16:AP:49:LEU:CD1	16:AP:51:VAL:HG23	2.30	0.61
25:AY:78:ALA:HA	25:AY:81:LYS:CD	2.31	0.61
27:B1:83:GLU:CG	27:B1:86:SER:N	2.63	0.61
35:BA:1411:C:HO2'	35:BA:1412:A:H8	1.47	0.61
35:BA:1625:C:H2'	35:BA:1626:G:O4'	2.01	0.61
35:BA:1823:G:O2'	35:BA:1824:G:H5'	2.00	0.61
35:BA:1854:A:H5'	35:BA:1855:G:OP2	2.00	0.61
35:BA:2775:A:O2'	35:BA:2776:A:H5'	1.99	0.61
35:BA:2830:G:H5'	39:BE:58:ARG:NH1	2.15	0.61
35:BA:92:A:H2'	35:BA:93:G:C8	2.35	0.61
42:BH:67:LEU:HG	42:BH:71:LEU:CD2	2.31	0.61
43:BI:3:VAL:HG12	43:BI:37:VAL:O	2.01	0.61
43:BI:71:ILE:HG13	43:BI:72:LEU:H	1.65	0.61
44:BN:120:LEU:C	44:BN:120:LEU:HD13	2.21	0.61
44:BN:56:ASN:CA	44:BN:124:ALA:HA	2.31	0.61
45:BO:42:SER:HA	45:BO:56:ASP:O	1.99	0.61
46:BP:64:LYS:O	46:BP:65:ARG:C	2.39	0.61
53:BW:35:ILE:HG22	53:BW:36:LEU:N	2.15	0.61
54:BX:77:LYS:CE	54:BX:78:LYS:HG3	2.30	0.61
54:BX:89:ILE:HA	54:BX:92:LEU:HD12	1.82	0.61
55:BY:75:ILE:HD11	55:BY:78:ALA:O	2.00	0.61
56:BZ:151:HIS:CD2	56:BZ:170:THR:HG22	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:552:U:O2'	1:CA:553:A:H5'	2.00	0.61
1:CA:685:G:N2	1:CA:686:U:H3	1.99	0.61
1:CA:715:A:H2'	1:CA:716:A:H8	1.64	0.61
1:CA:741:G:H2'	1:CA:742:G:O4'	2.01	0.61
5:CE:122:GLU:O	5:CE:123:LEU:HD23	2.01	0.61
7:CG:105:VAL:HG12	7:CG:109:ASN:ND2	2.16	0.61
12:CL:89:ARG:HE	12:CL:91:LYS:HE2	1.65	0.61
30:D4:29:PRO:C	30:D4:31:ILE:H	2.03	0.61
35:DA:1998:G:H2'	35:DA:1999:C:C6	2.36	0.61
26:D0:14:ARG:HD2	35:DA:2279:G:O6	2.00	0.61
35:DA:2348:U:H2'	35:DA:2349:G:C5'	2.26	0.61
35:DA:996:A:H2'	35:DA:997:G:H8	1.66	0.61
38:DD:94:LEU:HA	38:DD:104:TYR:HA	1.83	0.61
39:DE:120:TRP:CE2	39:DE:155:LYS:HB3	2.36	0.61
41:DG:39:ILE:CA	41:DG:157:ILE:HA	2.25	0.61
41:DG:171:ALA:O	41:DG:173:LEU:N	2.34	0.61
42:DH:122:THR:O	42:DH:133:VAL:HG13	2.01	0.61
42:DH:41:MET:CE	42:DH:55:PRO:HD2	2.29	0.61
44:DN:4:TYR:CD1	44:DN:4:TYR:N	2.68	0.61
47:DQ:53:ALA:HA	47:DQ:56:ARG:HB3	1.81	0.61
51:DU:96:ALA:O	51:DU:98:LEU:N	2.34	0.61
1:AA:383:A:H8	1:AA:383:A:O5'	1.84	0.60
1:AA:519:C:H2'	1:AA:520:A:C8	2.37	0.60
1:AA:547:A:H4'	1:AA:548:G:O5'	1.99	0.60
1:AA:640:A:H4'	8:AH:116:LYS:NZ	2.16	0.60
3:AC:11:ARG:HG2	3:AC:11:ARG:HH11	1.65	0.60
3:AC:182:ILE:HG12	3:AC:203:PHE:HA	1.83	0.60
4:AD:50:ARG:HD2	4:AD:51:PRO:O	2.01	0.60
7:AG:103:TRP:O	7:AG:107:ALA:N	2.17	0.60
9:AI:126:SER:O	9:AI:127:LYS:HB3	2.01	0.60
9:AI:79:LEU:O	9:AI:79:LEU:HD13	2.01	0.60
12:AL:26:ALA:O	12:AL:27:LEU:HB2	2.01	0.60
12:AL:55:VAL:HG12	12:AL:56:ALA:N	2.12	0.60
27:B1:15:ALA:O	27:B1:46:LEU:HD23	2.01	0.60
28:B2:32:LEU:C	28:B2:32:LEU:HD12	2.21	0.60
29:B3:6:VAL:HB	29:B3:54:VAL:HG11	1.83	0.60
31:B5:13:LYS:O	31:B5:16:ARG:HB3	2.01	0.60
35:BA:1031:G:N2	35:BA:1124:C:H1'	2.15	0.60
35:BA:2491:U:H4'	35:BA:2570:G:OP1	1.99	0.60
33:B7:5:TRP:CZ3	35:BA:464:U:H4'	2.36	0.60
26:B0:26:TYR:CE2	35:BA:857:C:H1'	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BB:65:C:C2'	36:BB:66:A:H5'	2.30	0.60
43:BI:99:GLU:O	43:BI:103:ARG:HB2	2.01	0.60
43:BI:94:ALA:C	43:BI:96:ASP:H	2.03	0.60
46:BP:50:ARG:CZ	46:BP:51:PHE:CZ	2.85	0.60
52:BV:24:LYS:HA	52:BV:94:LEU:HG	1.82	0.60
54:BX:39:ILE:HD12	54:BX:40:LYS:N	2.15	0.60
56:BZ:134:PRO:HG3	56:BZ:161:VAL:HG21	1.83	0.60
1:CA:110:C:H2'	1:CA:111:G:O4'	2.00	0.60
1:CA:1203:C:H2'	1:CA:1204:A:C8	2.36	0.60
1:CA:1218:C:H2'	1:CA:1219:U:C6	2.36	0.60
1:CA:1241:G:H2'	1:CA:1242:C:C5	2.36	0.60
1:CA:1323:G:H4'	1:CA:1363:C:C2	2.36	0.60
1:CA:1385:G:O2'	1:CA:1386:G:H5'	2.00	0.60
1:CA:865:A:H5'	1:CA:1078:U:C5	2.35	0.60
3:CC:182:ILE:HG23	3:CC:203:PHE:CA	2.32	0.60
6:CF:28:ARG:HG3	6:CF:28:ARG:HH11	1.65	0.60
7:CG:103:TRP:O	7:CG:107:ALA:N	2.17	0.60
23:CW:41:C:O2'	23:CW:42:C:H5'	2.01	0.60
25:CY:132:ILE:O	25:CY:136:ALA:N	2.32	0.60
25:CY:64:ARG:HA	25:CY:103:ILE:HD11	1.83	0.60
27:D1:19:GLN:CD	27:D1:44:PRO:HG3	2.20	0.60
27:D1:88:LYS:HA	27:D1:91:LYS:HB2	1.83	0.60
27:D1:89:GLU:CD	27:D1:89:GLU:H	2.03	0.60
28:D2:26:ARG:O	28:D2:29:LYS:HB2	2.01	0.60
29:D3:56:VAL:O	29:D3:57:GLU:HB2	2.01	0.60
31:D5:16:ARG:CG	31:D5:16:ARG:HH11	2.14	0.60
35:DA:1246:A:OP2	46:DP:16:ARG:NH2	2.34	0.60
35:DA:1997:G:O2'	35:DA:1998:G:H5'	2.01	0.60
35:DA:513:A:H1'	51:DU:11:ARG:NH1	2.16	0.60
35:DA:979:G:H3'	35:DA:980:A:C5'	2.31	0.60
38:DD:142:VAL:HG23	38:DD:193:VAL:N	2.15	0.60
38:DD:48:ARG:NH1	38:DD:48:ARG:HG3	2.16	0.60
40:DF:177:ALA:HB1	40:DF:178:PRO:CD	2.31	0.60
41:DG:2:PRO:O	41:DG:3:LEU:HB2	2.01	0.60
41:DG:47:LYS:O	41:DG:48:GLU:O	2.19	0.60
43:DI:31:LEU:HB3	43:DI:32:PRO:HD3	1.83	0.60
45:DO:68:GLU:OE2	45:DO:78:ARG:HD3	2.01	0.60
47:DQ:127:ILE:HD12	47:DQ:127:ILE:N	2.16	0.60
50:DT:74:ARG:HG2	50:DT:74:ARG:HH11	1.66	0.60
50:DT:86:ILE:HG12	50:DT:87:ASP:O	2.01	0.60
55:DY:2:ARG:C	55:DY:4:LYS:H	2.04	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1267:C:O2	1:AA:1267:C:H2'	2.01	0.60
1:AA:1277:C:C2'	1:AA:1278:U:H5'	2.31	0.60
1:AA:304:U:H2'	1:AA:305:G:C8	2.36	0.60
1:AA:1056:U:H5'	3:AC:163:ALA:CB	2.31	0.60
5:AE:11:ILE:HD12	5:AE:31:LEU:HD22	1.82	0.60
5:AE:150:ARG:CB	5:AE:150:ARG:HH11	2.14	0.60
6:AF:87:ARG:HH11	6:AF:87:ARG:HG3	1.65	0.60
25:AY:150:SER:C	25:AY:152:ASP:N	2.54	0.60
25:AY:65:THR:HG22	25:AY:66:LEU:N	2.16	0.60
27:B1:46:LEU:HD13	27:B1:46:LEU:O	2.00	0.60
27:B1:86:SER:CA	27:B1:89:GLU:CG	2.77	0.60
35:BA:2714:G:O2'	35:BA:2715:C:H5'	2.01	0.60
35:BA:708:C:N4	35:BA:723:G:H1	1.95	0.60
39:BE:142:GLY:C	39:BE:143:ASN:ND2	2.54	0.60
40:BF:117:ARG:HH21	40:BF:187:VAL:HA	1.65	0.60
41:BG:53:LEU:HD22	41:BG:53:LEU:N	2.14	0.60
42:BH:158:HIS:NE2	42:BH:170:ARG:HA	2.16	0.60
42:BH:89:ILE:CD1	42:BH:129:THR:HB	2.30	0.60
47:BQ:76:LYS:H	47:BQ:88:GLY:HA2	1.65	0.60
52:BV:64:HIS:CB	52:BV:96:ILE:HA	2.25	0.60
52:BV:19:LYS:HB3	52:BV:96:ILE:O	2.01	0.60
53:BW:5:ALA:HB3	53:BW:105:VAL:H	1.66	0.60
54:BX:64:LYS:CG	54:BX:65:ARG:N	2.64	0.60
54:BX:70:LEU:HG	54:BX:71:GLY:N	2.15	0.60
1:CA:1106:G:H2'	1:CA:1107:C:H6	1.65	0.60
1:CA:1363(A):A:H4'	1:CA:1364:U:C5'	2.26	0.60
1:CA:186:C:H4'	20:CT:82:SER:HB3	1.82	0.60
1:CA:262:A:H2'	1:CA:263:A:C8	2.36	0.60
1:CA:291:C:O2'	1:CA:292:G:H5'	2.02	0.60
1:CA:414:A:H2'	1:CA:415:A:O4'	2.02	0.60
1:CA:449:C:H2'	1:CA:450:G:O4'	2.01	0.60
1:CA:628:G:O2'	1:CA:629:G:H5'	2.01	0.60
1:CA:710:G:H5''	6:CF:54:LYS:HE3	1.81	0.60
1:CA:831:U:H2'	1:CA:832:C:C6	2.36	0.60
2:CB:102:LEU:HD12	2:CB:102:LEU:H	1.65	0.60
2:CB:185:ILE:HG22	2:CB:199:TYR:CD1	2.36	0.60
3:CC:206:GLU:O	3:CC:208:ILE:N	2.34	0.60
9:CI:114:TYR:CE1	10:CJ:59:SER:HA	2.35	0.60
1:CA:473:G:H5''	16:CP:81:ARG:HE	1.66	0.60
25:CY:7:TYR:CZ	25:CY:160:GLU:HG2	2.36	0.60
35:DA:528:A:C2	35:DA:2043:C:C5'	2.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2248:C:C2'	35:DA:2249:U:H5'	2.31	0.60
35:DA:2624:G:O2'	35:DA:2625:G:H5'	2.01	0.60
35:DA:2708:G:H2'	35:DA:2709:G:H8	1.65	0.60
35:DA:271(E):U:H2'	35:DA:271(F):C:C6	2.36	0.60
35:DA:260:G:H1'	35:DA:621:A:H1'	1.83	0.60
38:DD:32:SER:OG	38:DD:33:LEU:N	2.34	0.60
38:DD:80:ALA:HB3	38:DD:94:LEU:HD13	1.83	0.60
41:DG:103:LEU:HB3	41:DG:107:LEU:HD12	1.82	0.60
41:DG:125:PHE:HB3	41:DG:128:ARG:O	2.01	0.60
44:DN:73:THR:O	44:DN:75:TYR:N	2.30	0.60
47:DQ:82:ARG:HG2	47:DQ:82:ARG:NH1	2.16	0.60
52:DV:64:HIS:CB	52:DV:96:ILE:HA	2.24	0.60
55:DY:37:VAL:HG22	55:DY:67:LEU:O	2.00	0.60
56:DZ:39:VAL:HG23	56:DZ:40:ASP:N	2.17	0.60
1:AA:1437:C:N4	1:AA:1464:G:H1	1.97	0.60
1:AA:501:C:H2'	1:AA:502:G:H8	1.66	0.60
1:AA:693:G:H2'	1:AA:694:A:C8	2.36	0.60
1:AA:831:U:H2'	1:AA:832:C:C6	2.36	0.60
2:AB:75:LYS:CA	2:AB:78:GLN:HE21	2.02	0.60
3:AC:141:VAL:O	3:AC:141:VAL:HG12	2.01	0.60
5:AE:147:ASP:HA	5:AE:150:ARG:NH1	2.11	0.60
8:AH:88:LYS:HB3	8:AH:89:PRO:HD2	1.83	0.60
17:AQ:76:LEU:HD12	17:AQ:77:VAL:H	1.66	0.60
19:AS:16:LEU:H	19:AS:16:LEU:CD1	2.14	0.60
25:AY:30:THR:HG21	25:AY:179:LYS:HD2	1.81	0.60
27:B1:48:LYS:CG	27:B1:49:VAL:H	2.08	0.60
27:B1:68:PRO:CG	27:B1:69:LYS:H	2.07	0.60
27:B1:68:PRO:O	27:B1:71:TYR:N	2.35	0.60
35:BA:1424:G:H2'	35:BA:1425:G:O4'	2.01	0.60
35:BA:2264:C:H2'	35:BA:2265:U:C6	2.36	0.60
35:BA:250:G:H2'	35:BA:251:A:C8	2.36	0.60
35:BA:2708:G:H2'	35:BA:2709:G:H8	1.66	0.60
35:BA:347:A:H2'	35:BA:348:G:C8	2.36	0.60
35:BA:736:C:H2'	35:BA:737:C:H6	1.67	0.60
36:BB:75:G:N2	56:BZ:87:ASP:OD2	2.35	0.60
38:BD:35:LYS:NZ	38:BD:104:TYR:HB2	2.16	0.60
38:BD:242:ARG:HG3	38:BD:242:ARG:NH1	2.14	0.60
40:BF:29:ASN:HB3	40:BF:112:MET:HE3	1.83	0.60
44:BN:73:THR:O	44:BN:75:TYR:N	2.28	0.60
45:BO:46:ALA:N	45:BO:54:GLU:HG2	2.17	0.60
48:BR:9:LYS:HZ1	48:BR:42:LYS:HB3	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:58:ASN:N	50:BT:58:ASN:HD22	1.98	0.60
51:BU:33:ARG:HA	51:BU:36:ARG:HB2	1.82	0.60
52:BV:1:MET:HE1	52:BV:46:VAL:HG23	1.83	0.60
55:BY:28:LYS:CE	55:BY:30:VAL:HA	2.31	0.60
56:BZ:120:ILE:HB	56:BZ:171:ILE:C	2.21	0.60
1:CA:1030(D):A:H2'	1:CA:1031:G:H5'	1.84	0.60
1:CA:1035:A:H2'	1:CA:1036:G:H8	1.66	0.60
1:CA:597:G:C2'	1:CA:598:U:H5'	2.30	0.60
2:CB:102:LEU:O	2:CB:105:PHE:HB2	2.01	0.60
4:CD:62:GLN:HB3	4:CD:66:ARG:NH2	2.16	0.60
8:CH:82:HIS:CD2	8:CH:138:TRP:NE1	2.68	0.60
9:CI:71:SER:HA	9:CI:74:ILE:HD12	1.82	0.60
11:CK:43:SER:CA	11:CK:47:VAL:HG21	2.31	0.60
31:D5:40:LYS:HE2	31:D5:46:CYS:CB	2.30	0.60
32:D6:13:CYS:HA	32:D6:50:ARG:O	2.01	0.60
35:DA:2208:A:H1'	35:DA:2219:G:C2	2.36	0.60
35:DA:2744:G:N7	35:DA:2755:C:O2	2.35	0.60
35:DA:2821:A:H3'	35:DA:2821:A:OP2	2.01	0.60
35:DA:778:G:H5''	38:DD:48:ARG:HD2	1.83	0.60
36:DB:42:C:H4'	41:DG:67:LYS:HB3	1.84	0.60
48:DR:52:ILE:O	48:DR:55:ALA:N	2.35	0.60
50:DT:118:ARG:O	50:DT:121:ILE:HG22	2.00	0.60
52:DV:79:VAL:HG12	52:DV:80:GLN:N	2.16	0.60
54:DX:29:TRP:CZ3	54:DX:76:ARG:HG2	2.34	0.60
28:D2:23:LYS:CA	54:DX:5:TYR:CE1	2.85	0.60
54:DX:64:LYS:CG	54:DX:65:ARG:N	2.64	0.60
54:DX:8:ILE:N	54:DX:8:ILE:HD12	2.15	0.60
55:DY:47:LYS:HG3	55:DY:60:PHE:CZ	2.36	0.60
55:DY:75:ILE:HD11	55:DY:78:ALA:O	2.01	0.60
1:AA:445:G:H2'	1:AA:446:G:H8	1.66	0.60
1:AA:501:C:O2'	1:AA:502:G:H5'	2.02	0.60
1:AA:627:G:H2'	1:AA:628:G:C8	2.35	0.60
2:AB:12:GLU:OE2	2:AB:214:ILE:HD11	2.00	0.60
2:AB:213:LEU:C	2:AB:213:LEU:HD23	2.22	0.60
3:AC:124:ILE:HG13	3:AC:130:VAL:HG22	1.83	0.60
4:AD:56:VAL:O	4:AD:58:LEU:N	2.34	0.60
4:AD:59:ARG:NH1	4:AD:59:ARG:HA	2.16	0.60
8:AH:122:ARG:HA	8:AH:125:ARG:CB	2.31	0.60
8:AH:26:VAL:HG22	8:AH:32:LYS:HZ2	1.65	0.60
9:AI:118:LYS:NZ	9:AI:118:LYS:HB3	2.15	0.60
11:AK:44:SER:O	11:AK:47:VAL:HB	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:22:SER:C	12:AL:24:VAL:H	2.04	0.60
1:AA:390:C:O3'	16:AP:28:ARG:NH2	2.34	0.60
32:B6:13:CYS:HA	32:B6:50:ARG:O	2.01	0.60
35:BA:1127:A:C2'	35:BA:1128:A:H5''	2.32	0.60
35:BA:1448:G:H2'	35:BA:1449:A:C8	2.37	0.60
35:BA:2175:C:H2'	35:BA:2176:A:C5'	2.29	0.60
35:BA:257:A:H2'	35:BA:258:G:O4'	2.00	0.60
38:BD:204:ILE:HG13	38:BD:204:ILE:O	2.00	0.60
38:BD:211:ARG:HH11	38:BD:211:ARG:CG	2.13	0.60
38:BD:227:ASN:HB3	38:BD:228:PRO:HD2	1.82	0.60
38:BD:57:GLY:HA2	38:BD:214:TRP:O	2.01	0.60
39:BE:104:VAL:O	39:BE:167:VAL:HG12	2.01	0.60
43:BI:94:ALA:HA	43:BI:97:ILE:CG1	2.31	0.60
46:BP:39:LYS:C	46:BP:41:ARG:H	2.04	0.60
50:BT:14:TYR:CD1	50:BT:14:TYR:N	2.68	0.60
50:BT:28:VAL:CG2	50:BT:47:GLY:H	2.07	0.60
35:BA:565:C:O3'	52:BV:81:TYR:HE1	1.84	0.60
55:BY:81:LYS:HG2	55:BY:97:ARG:H	1.66	0.60
1:CA:460:G:O6	1:CA:470:C:H5''	2.02	0.60
4:CD:79:PHE:HA	4:CD:93:PHE:CD2	2.37	0.60
15:CO:11:VAL:HG21	15:CO:34:LEU:HD23	1.82	0.60
17:CQ:95:TYR:C	17:CQ:97:SER:H	2.03	0.60
23:CW:59:A:C2	23:CW:61:U:H2'	2.37	0.60
27:D1:16:ASN:O	27:D1:16:ASN:ND2	2.34	0.60
35:DA:347:A:H2'	35:DA:348:G:C8	2.36	0.60
35:DA:618:C:H2'	35:DA:619:G:O4'	2.01	0.60
35:DA:853:G:H2'	35:DA:854:G:C8	2.36	0.60
36:DB:15:A:H3'	36:DB:16:G:H5'	1.83	0.60
41:DG:82:LEU:HB3	41:DG:87:PRO:HG3	1.83	0.60
42:DH:146:ALA:O	42:DH:147:ASN:C	2.38	0.60
46:DP:7:ARG:HB3	46:DP:8:PRO:CD	2.30	0.60
47:DQ:61:GLY:H	56:DZ:177:PRO:HB2	1.66	0.60
1:CA:1442(A):G:H22	50:DT:119:LYS:HB2	1.65	0.60
50:DT:11:GLU:CD	50:DT:11:GLU:N	2.39	0.60
56:DZ:8:TYR:HA	56:DZ:62:PRO:HG2	1.83	0.60
1:AA:625:G:C4	1:AA:626:U:C5	2.90	0.60
1:AA:834:C:H2'	1:AA:835:U:H6	1.66	0.60
2:AB:169:LYS:HD3	2:AB:169:LYS:O	2.01	0.60
2:AB:31:TYR:HD2	2:AB:31:TYR:N	2.00	0.60
8:AH:26:VAL:O	8:AH:59:LEU:N	2.34	0.60
10:AJ:40:LEU:HD12	10:AJ:41:PRO:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:36:VAL:O	12:AL:58:VAL:HG13	2.02	0.60
16:AP:20:VAL:HG23	16:AP:34:GLU:O	2.01	0.60
35:BA:1847:A:H2'	35:BA:1847:A:N3	2.16	0.60
27:B1:40:ARG:HD2	35:BA:2081:C:H4'	1.84	0.60
35:BA:2626:C:H2'	35:BA:2627:G:H8	1.65	0.60
35:BA:393:C:H2'	35:BA:394:A:H8	1.66	0.60
35:BA:632:A:H2'	35:BA:633:A:C8	2.37	0.60
35:BA:869:G:H2'	35:BA:870:A:H8	1.66	0.60
38:BD:222:ARG:O	38:BD:224:ALA:N	2.35	0.60
35:BA:2632:A:O2'	39:BE:61:ARG:NH2	2.35	0.60
42:BH:89:ILE:N	42:BH:89:ILE:CD1	2.64	0.60
47:BQ:127:ILE:N	47:BQ:127:ILE:HD12	2.15	0.60
48:BR:12:ARG:HG3	48:BR:12:ARG:HH11	1.66	0.60
49:BS:28:VAL:H	49:BS:89:ARG:HB2	1.66	0.60
56:BZ:105:VAL:N	56:BZ:141:VAL:HG11	2.17	0.60
56:BZ:41:LEU:O	56:BZ:44:PHE:HB3	2.01	0.60
1:CA:1343:G:H1'	9:CI:121:ARG:HH12	1.66	0.60
1:CA:148:G:H1	1:CA:174:C:H42	1.49	0.60
1:CA:460:G:H21	1:CA:472:A:H62	1.49	0.60
2:CB:107:THR:HG23	2:CB:110:GLN:OE1	2.02	0.60
4:CD:106:TYR:CE1	4:CD:113:SER:HA	2.37	0.60
4:CD:176:LEU:HG	4:CD:177:ASP:N	2.16	0.60
5:CE:78:HIS:HB2	5:CE:79:GLU:OE1	2.01	0.60
9:CI:10:ARG:NH2	9:CI:107:ARG:HD3	2.16	0.60
9:CI:26:VAL:HG22	9:CI:61:ALA:HB3	1.84	0.60
10:CJ:49:VAL:HG13	14:CN:41:ARG:HB2	1.82	0.60
15:CO:8:LYS:O	15:CO:12:ILE:HG13	2.01	0.60
25:CY:30:THR:C	25:CY:32:ARG:H	2.04	0.60
35:DA:1192:G:C2'	35:DA:1193:G:H5'	2.32	0.60
35:DA:1688:U:H5'	35:DA:1689:A:OP1	2.02	0.60
35:DA:1811:G:O2'	35:DA:1812:A:H5'	2.02	0.60
35:DA:2555:U:H2'	35:DA:2556:C:H5'	1.82	0.60
35:DA:271(J):C:H2'	35:DA:271(K):U:H5''	1.83	0.60
35:DA:542:C:H2'	35:DA:543:C:OP1	2.02	0.60
35:DA:64:A:H2'	35:DA:65:C:C6	2.36	0.60
35:DA:998:C:P	51:DU:93:LYS:HE2	2.41	0.60
37:DC:82:LYS:HB3	37:DC:86:ALA:HB2	1.83	0.60
42:DH:144:VAL:HG12	42:DH:148:ILE:HD11	1.83	0.60
42:DH:67:LEU:HG	42:DH:71:LEU:CD2	2.30	0.60
44:DN:42:TRP:CE2	44:DN:44:PRO:HD3	2.36	0.60
49:DS:85:VAL:CG2	49:DS:106:ARG:HB2	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1010:G:H22	1:AA:1020:U:H1'	1.65	0.60
1:AA:1488:G:H2'	1:AA:1489:G:H8	1.67	0.60
1:AA:519:C:H2'	1:AA:520:A:H8	1.65	0.60
1:AA:734:G:O2'	1:AA:735:C:H5'	2.01	0.60
1:AA:811:C:H4'	1:AA:900:A:N6	2.17	0.60
1:AA:9:G:H5'	5:AE:122:GLU:OE2	2.01	0.60
2:AB:79:ASP:O	2:AB:82:ARG:N	2.34	0.60
4:AD:170:VAL:HG13	4:AD:171:GLY:N	2.16	0.60
4:AD:176:LEU:CD1	4:AD:177:ASP:H	2.14	0.60
8:AH:51:VAL:HG11	8:AH:60:ARG:HG2	1.81	0.60
13:AM:108:ARG:N	13:AM:108:ARG:HD2	2.03	0.60
13:AM:13:LYS:HZ2	13:AM:21:TYR:HE1	1.49	0.60
27:B1:88:LYS:HA	27:B1:91:LYS:HD2	1.84	0.60
29:B3:56:VAL:CG1	29:B3:57:GLU:H	2.00	0.60
35:BA:1642:G:H2'	35:BA:1643:G:H8	1.65	0.60
35:BA:1703:G:H2'	35:BA:1704:G:C8	2.37	0.60
35:BA:2339:G:H2'	35:BA:2340:G:H8	1.67	0.60
35:BA:37:C:H2'	35:BA:38:A:C8	2.35	0.60
37:BC:44:HIS:HA	37:BC:174:PRO:CB	2.31	0.60
38:BD:206:LEU:HD23	38:BD:211:ARG:HH11	1.66	0.60
40:BF:89:VAL:HG12	40:BF:90:PHE:N	2.16	0.60
42:BH:102:ALA:HB1	42:BH:115:VAL:O	2.02	0.60
42:BH:46:GLU:O	42:BH:47:GLU:HB2	2.02	0.60
35:BA:2748:A:H2	42:BH:63:SER:HB3	1.67	0.60
43:BI:91:SER:CB	43:BI:121:LYS:HE3	2.30	0.60
44:BN:58:ASP:OD1	44:BN:124:ALA:HB1	2.01	0.60
44:BN:66:LYS:O	44:BN:87:LEU:HB3	2.02	0.60
45:BO:68:GLU:OE2	45:BO:78:ARG:HD3	2.01	0.60
46:BP:108:LYS:O	46:BP:110:TYR:N	2.31	0.60
46:BP:39:LYS:HD3	46:BP:40:SER:N	2.16	0.60
47:BQ:43:THR:HG1	47:BQ:46:GLN:HG3	1.65	0.60
51:BU:90:VAL:CG2	52:BV:39:LEU:HD12	2.32	0.60
54:BX:82:GLN:HB3	54:BX:85:PRO:HG2	1.83	0.60
56:BZ:6:LYS:HA	56:BZ:60:GLU:O	2.01	0.60
1:CA:105:G:H2'	1:CA:106:C:C6	2.36	0.60
4:CD:50:ARG:HD2	4:CD:51:PRO:O	2.00	0.60
5:CE:150:ARG:CB	5:CE:150:ARG:HH11	2.14	0.60
9:CI:118:LYS:HB3	9:CI:118:LYS:NZ	2.16	0.60
12:CL:24:VAL:HG13	12:CL:98:TYR:HE2	1.63	0.60
16:CP:6:LEU:N	16:CP:6:LEU:CD1	2.65	0.60
18:CR:70:ILE:O	18:CR:74:ARG:HB2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:CY:88:LEU:H	25:CY:88:LEU:HD12	1.65	0.60
27:D1:23:LYS:HB3	27:D1:37:ILE:CD1	2.24	0.60
31:D5:12:SER:O	31:D5:13:LYS:C	2.38	0.60
31:D5:30:LEU:HD11	53:DW:38:TYR:HB2	1.83	0.60
34:D8:36:LYS:O	34:D8:37:SER:O	2.19	0.60
34:D8:43:GLN:O	34:D8:44:LYS:HD2	2.02	0.60
35:DA:1021:A:H3'	35:DA:1021:A:H8	1.67	0.60
35:DA:143:G:H2'	35:DA:143(A):C:C6	2.36	0.60
35:DA:143(A):C:H2'	35:DA:143(A):C:O2	2.01	0.60
35:DA:2103:C:C2'	35:DA:2104:G:H5''	2.31	0.60
35:DA:2262:U:H4'	35:DA:2328:A:C2	2.36	0.60
35:DA:2328:A:H2'	35:DA:2329:G:C8	2.36	0.60
35:DA:648:G:C4'	35:DA:2351:G:H5''	2.32	0.60
35:DA:244:A:H2'	35:DA:245:G:O4'	2.01	0.60
35:DA:2742:C:O2'	35:DA:2743:C:H5'	2.02	0.60
35:DA:2745:C:H2'	35:DA:2746:U:H6	1.65	0.60
35:DA:2759:G:O2'	35:DA:2760:C:H5'	2.02	0.60
35:DA:2876:G:H4'	50:DT:2:ASN:O	2.02	0.60
35:DA:792:G:C4'	35:DA:793:A:H5'	2.31	0.60
35:DA:878:A:H3'	35:DA:879:G:H8	1.65	0.60
37:DC:59:ARG:HH21	37:DC:199:HIS:CB	2.15	0.60
37:DC:76:ALA:HB3	37:DC:94:VAL:HG11	1.84	0.60
39:DE:59:VAL:HG21	39:DE:63:LEU:HA	1.84	0.60
40:DF:128:ALA:O	40:DF:142:TRP:NE1	2.33	0.60
41:DG:158:ALA:O	41:DG:159:VAL:HG13	2.01	0.60
41:DG:96:ARG:HG3	41:DG:97:ASP:N	2.16	0.60
43:DI:14:ASP:O	43:DI:15:VAL:O	2.19	0.60
44:DN:3:THR:O	44:DN:5:VAL:N	2.27	0.60
44:DN:44:PRO:C	44:DN:46:VAL:H	2.04	0.60
48:DR:29:LEU:HD23	48:DR:70:LEU:HD11	1.84	0.60
48:DR:6:SER:HA	48:DR:8:ARG:NH2	2.17	0.60
35:DA:1252:G:N3	51:DU:33:ARG:HD2	2.17	0.60
52:DV:22:VAL:HG21	52:DV:96:ILE:HD12	1.84	0.60
54:DX:28:PHE:CD1	54:DX:28:PHE:N	2.69	0.60
56:DZ:86:VAL:HG12	56:DZ:87:ASP:H	1.65	0.60
1:AA:1281:U:H4'	1:AA:1282:C:OP2	2.01	0.60
1:AA:1313:U:P	19:AS:6:LYS:HG3	2.42	0.60
1:AA:1323:G:H4'	1:AA:1363:C:C2	2.36	0.60
1:AA:688:G:H5'	11:AK:46:GLY:O	2.02	0.60
2:AB:200:ILE:O	2:AB:201:ILE:HD13	2.02	0.60
5:AE:91:LEU:HG	5:AE:120:THR:HG22	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:26:ILE:O	6:AF:30:LEU:HG	2.01	0.60
15:AO:23:GLY:O	15:AO:24:SER:HB3	2.01	0.60
31:B5:32:PRO:O	31:B5:33:CYS:HB2	2.00	0.60
31:B5:52:TYR:HA	31:B5:56:LYS:HZ2	1.67	0.60
34:B8:49:VAL:HB	34:B8:53:PRO:HD3	1.82	0.60
35:BA:1232:G:H2'	35:BA:1233:C:C6	2.36	0.60
35:BA:1615:C:H5	35:BA:1617:C:C4	2.19	0.60
35:BA:271(E):U:H2'	35:BA:271(F):C:C6	2.37	0.60
35:BA:2825:C:H2'	35:BA:2826:A:O4'	2.02	0.60
35:BA:838:C:H42	35:BA:940:G:H1	1.49	0.60
37:BC:45:ALA:H	37:BC:174:PRO:CB	2.14	0.60
40:BF:158:THR:HG21	40:BF:163:VAL:CB	2.26	0.60
40:BF:177:ALA:HB1	40:BF:178:PRO:CD	2.30	0.60
35:BA:673:C:H5'	40:BF:54:ARG:NH1	2.17	0.60
43:BI:133:HIS:CB	43:BI:134:PRO:CD	2.78	0.60
44:BN:34:LEU:HD23	44:BN:120:LEU:HD23	1.84	0.60
45:BO:31:LYS:C	45:BO:32:TYR:CD1	2.73	0.60
1:AA:1422:G:H4'	45:BO:49:ARG:NH1	2.17	0.60
49:BS:104:GLY:O	49:BS:106:ARG:N	2.28	0.60
49:BS:41:ASP:O	49:BS:45:GLY:HA2	2.00	0.60
51:BU:18:LEU:CD2	51:BU:22:LYS:HE2	2.31	0.60
53:BW:9:TYR:HD2	53:BW:102:HIS:HE2	1.49	0.60
54:BX:88:LYS:HD2	54:BX:88:LYS:N	2.17	0.60
1:CA:119:A:O2'	1:CA:120:A:OP2	2.19	0.60
1:CA:1513:A:C6	1:CA:1514:C:N4	2.69	0.60
1:CA:490:G:H2'	1:CA:491:G:C8	2.36	0.60
1:CA:775:G:O2'	1:CA:776:G:H5'	2.02	0.60
1:CA:884:U:H4'	1:CA:885:G:C5'	2.32	0.60
4:CD:18:LYS:NZ	4:CD:31:CYS:HB3	2.16	0.60
10:CJ:29:ARG:HH11	10:CJ:29:ARG:HG2	1.66	0.60
10:CJ:40:LEU:HD12	10:CJ:41:PRO:O	2.01	0.60
10:CJ:35:SER:O	10:CJ:72:VAL:HG13	2.01	0.60
1:CA:1227:A:OP2	13:CM:111:LYS:HE3	2.00	0.60
25:CY:128:ALA:O	25:CY:132:ILE:HG13	2.01	0.60
29:D3:56:VAL:O	29:D3:57:GLU:CB	2.50	0.60
35:DA:151:C:H42	35:DA:175:G:H1	1.50	0.60
35:DA:1765:C:H2'	35:DA:1766:U:C6	2.36	0.60
35:DA:2875:C:H4'	50:DT:5:ALA:HB2	1.82	0.60
35:DA:291:C:H2'	35:DA:292:C:H6	1.64	0.60
35:DA:622:G:O2'	35:DA:623:G:H5'	2.02	0.60
36:DB:104:U:O2'	56:DZ:72:ARG:HD2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:130:ALA:HB2	38:DD:192:THR:CB	2.32	0.60
40:DF:141:ALA:O	40:DF:144:LYS:HB3	2.01	0.60
41:DG:104:GLU:O	41:DG:106:LEU:N	2.34	0.60
42:DH:68:THR:HA	42:DH:71:LEU:HD22	1.83	0.60
42:DH:87:LEU:C	42:DH:88:LEU:HD22	2.22	0.60
45:DO:3:GLN:HB2	45:DO:4:PRO:HD2	1.84	0.60
46:DP:112:LEU:C	46:DP:112:LEU:HD13	2.21	0.60
50:DT:14:TYR:HD1	50:DT:14:TYR:N	1.99	0.60
51:DU:52:ARG:O	51:DU:54:LYS:N	2.35	0.60
52:DV:96:ILE:CG2	52:DV:97:LYS:N	2.64	0.60
54:DX:36:LYS:C	54:DX:38:GLU:H	2.04	0.60
56:DZ:118:GLN:NE2	56:DZ:118:GLN:HA	2.17	0.60
1:AA:1294:G:H2'	1:AA:1295:G:H8	1.67	0.60
1:AA:1104:G:OP1	2:AB:111:ARG:HD2	2.01	0.60
2:AB:185:ILE:HG22	2:AB:199:TYR:HD1	1.67	0.60
2:AB:168:THR:HG23	2:AB:192:SER:OG	2.01	0.60
2:AB:224:GLN:HG2	2:AB:224:GLN:O	2.02	0.60
4:AD:108:LEU:CB	4:AD:110:PHE:CE1	2.84	0.60
4:AD:9:CYS:SG	4:AD:31:CYS:O	2.59	0.60
5:AE:69:VAL:O	5:AE:71:LEU:N	2.35	0.60
7:AG:86:GLN:HG2	23:AW:33:C:C5'	2.32	0.60
9:AI:113:LYS:N	9:AI:113:LYS:HD2	2.16	0.60
10:AJ:50:ILE:HA	10:AJ:60:ARG:HG2	1.82	0.60
12:AL:115:LYS:O	12:AL:117:ARG:HG3	2.01	0.60
15:AO:33:THR:OG1	15:AO:63:ARG:HD2	2.02	0.60
15:AO:82:ILE:HD11	15:AO:87:ILE:O	2.02	0.60
17:AQ:5:VAL:HA	17:AQ:59:ILE:O	2.02	0.60
22:AV:30:A:H2'	22:AV:31:U:C5	2.37	0.60
25:AY:68:VAL:O	25:AY:70:SER:N	2.33	0.60
31:B5:52:TYR:HA	31:B5:56:LYS:NZ	2.17	0.60
35:BA:1280:G:C3'	35:BA:1281:G:H5''	2.32	0.60
35:BA:1440:G:H2'	35:BA:1441:G:C8	2.36	0.60
35:BA:2121:G:H1	35:BA:2177:C:H42	1.49	0.60
35:BA:2223:G:H2'	35:BA:2224:G:H5'	1.84	0.60
35:BA:244:A:H2'	35:BA:245:G:O4'	2.02	0.60
35:BA:2820:A:O3'	48:BR:2:ARG:NH2	2.34	0.60
35:BA:861:A:H62	35:BA:916:G:H21	1.48	0.60
38:BD:231:HIS:ND1	38:BD:232:PRO:CD	2.64	0.60
38:BD:97:TYR:HB2	38:BD:101:GLU:O	2.01	0.60
40:BF:178:PRO:CG	40:BF:179:GLU:H	2.12	0.60
40:BF:3:GLU:CB	40:BF:24:LEU:HG	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BQ:68:ILE:N	47:BQ:68:ILE:HD13	2.14	0.60
1:CA:1277:C:C2'	1:CA:1278:U:H5'	2.32	0.60
1:CA:393:A:OP2	16:CP:12:LYS:HD3	2.01	0.60
1:CA:818:G:C3'	1:CA:819:A:H5''	2.31	0.60
1:CA:923:A:H2'	1:CA:924:C:C6	2.36	0.60
2:CB:28:PHE:CE1	2:CB:31:TYR:HB2	2.36	0.60
8:CH:18:ARG:N	8:CH:78:GLN:HE22	1.99	0.60
1:CA:1251:A:H5''	9:CI:12:GLU:OE1	2.01	0.60
10:CJ:32:ALA:HB1	10:CJ:75:ILE:CD1	2.31	0.60
11:CK:99:GLN:C	11:CK:101:SER:H	2.05	0.60
12:CL:46:LYS:CG	12:CL:47:LYS:H	2.09	0.60
1:CA:974:A:C1'	14:CN:31:ARG:HH21	2.15	0.60
14:CN:40:CYS:SG	14:CN:41:ARG:N	2.74	0.60
16:CP:21:VAL:O	16:CP:33:ILE:HB	2.02	0.60
18:CR:36:ASN:HD22	18:CR:39:VAL:CB	2.14	0.60
25:CY:18:LEU:O	25:CY:20:VAL:N	2.34	0.60
31:D5:40:LYS:HZ2	31:D5:45:VAL:HA	1.67	0.60
35:DA:1177:A:H5''	35:DA:1178:C:O5'	2.01	0.60
35:DA:122:G:H1	35:DA:129:C:H42	1.50	0.60
35:DA:1467:C:OP2	35:DA:1547:C:H5	1.85	0.60
35:DA:1689:A:N6	35:DA:1698:A:H2	1.99	0.60
35:DA:1847:A:N3	35:DA:1847:A:H2'	2.16	0.60
35:DA:2179:C:H1'	37:DC:171:ILE:CB	2.32	0.60
35:DA:2282:G:H1	35:DA:2427:C:N4	1.94	0.60
35:DA:2314:C:H2'	35:DA:2315:G:H8	1.67	0.60
35:DA:2766:G:N3	35:DA:2766:G:H2'	2.16	0.60
35:DA:2864:G:O2'	35:DA:2865:U:H5'	2.02	0.60
33:D7:5:TRP:CZ3	35:DA:464:U:H4'	2.37	0.60
38:DD:10:THR:O	38:DD:11:PRO:C	2.35	0.60
38:DD:222:ARG:O	38:DD:223:GLY:C	2.40	0.60
41:DG:96:ARG:O	41:DG:99:MET:HB3	2.01	0.60
42:DH:89:ILE:HD13	42:DH:90:LYS:H	1.66	0.60
54:DX:35:THR:HB	54:DX:75:ASP:OD2	2.02	0.60
54:DX:89:ILE:HD12	54:DX:89:ILE:N	2.16	0.60
1:AA:532:A:H2	1:AA:1207:G:O4'	1.84	0.60
1:AA:1392:G:N2	1:AA:1502:A:H8	2.00	0.60
1:AA:711:G:O2'	1:AA:712:A:H5'	2.02	0.60
1:AA:865:A:H2'	1:AA:866:C:O4'	2.02	0.60
2:AB:172:ILE:HD12	2:AB:172:ILE:N	2.10	0.60
3:AC:138:VAL:CG2	3:AC:151:VAL:HG23	2.32	0.60
6:AF:69:GLU:HG2	6:AF:70:ASP:H	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:121:ALA:O	7:AG:124:LEU:HB2	2.02	0.60
7:AG:21:VAL:HG23	7:AG:22:LEU:H	1.66	0.60
16:AP:53:VAL:HG23	16:AP:54:GLU:H	1.67	0.60
28:B2:12:GLU:C	28:B2:14:ARG:HE	2.04	0.60
35:BA:1238:G:H2'	35:BA:1239:G:H8	1.67	0.60
35:BA:141:A:C8	35:BA:1408:C:O2'	2.54	0.60
35:BA:1555:G:H2'	35:BA:1556:C:C6	2.37	0.60
35:BA:1720:U:C2'	35:BA:1721:G:H5'	2.32	0.60
35:BA:2742:C:O2'	35:BA:2743:C:H5'	2.01	0.60
35:BA:2627:G:N3	35:BA:2781:A:H2	1.99	0.60
35:BA:291:C:H2'	35:BA:292:C:H6	1.66	0.60
35:BA:523:C:O2'	35:BA:524:U:H5'	2.02	0.60
35:BA:674:G:P	40:BF:54:ARG:HH22	2.25	0.60
35:BA:692:C:O2'	35:BA:693:C:H5'	2.02	0.60
44:BN:55:VAL:HG12	44:BN:126:PRO:HA	1.84	0.60
46:BP:13:ASN:HD22	46:BP:13:ASN:H	1.50	0.60
47:BQ:69:PHE:CD1	47:BQ:70:PRO:HD2	2.36	0.60
54:BX:36:LYS:C	54:BX:38:GLU:H	2.04	0.60
1:CA:967:C:H2'	1:CA:968:A:N7	2.17	0.60
2:CB:168:THR:HG21	2:CB:191:ASP:OD1	2.01	0.60
5:CE:76:ILE:HG23	5:CE:77:PRO:N	2.17	0.60
1:CA:710:G:H5''	6:CF:54:LYS:CE	2.32	0.60
1:CA:255:G:O3'	17:CQ:17:LYS:HD2	2.02	0.60
6:CF:62:TRP:CB	18:CR:35:ARG:HH12	2.15	0.60
25:CY:169:ILE:O	25:CY:172:ALA:N	2.35	0.60
25:CY:36:ALA:HA	25:CY:39:LEU:CG	2.31	0.60
27:D1:48:LYS:O	27:D1:49:VAL:HG23	2.01	0.60
29:D3:1:MET:SD	29:D3:38:GLU:HG2	2.42	0.60
29:D3:6:VAL:HB	29:D3:54:VAL:HG11	1.83	0.60
33:D7:30:VAL:HG23	33:D7:31:LEU:N	2.15	0.60
35:DA:1041:C:H5'	35:DA:1042:G:OP2	2.02	0.60
35:DA:143:G:H2'	35:DA:143(A):C:H6	1.67	0.60
35:DA:1602:U:H3'	35:DA:1603:A:H5''	1.83	0.60
35:DA:1909:C:O2	35:DA:1909:C:H2'	2.00	0.60
35:DA:2600:A:C2'	35:DA:2601:C:H5'	2.32	0.60
33:D7:16:HIS:CE1	35:DA:465:G:H4'	2.37	0.60
38:DD:271:ILE:N	38:DD:271:ILE:HD12	2.17	0.60
42:DH:137:ASP:O	42:DH:138:LYS:HB2	2.02	0.60
45:DO:36:GLY:H	45:DO:62:VAL:HB	1.67	0.60
46:DP:108:LYS:C	46:DP:110:TYR:H	2.05	0.60
46:DP:99:LEU:O	46:DP:102:ARG:HB3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:28:VAL:H	49:DS:89:ARG:HB2	1.66	0.60
50:DT:109:GLU:CA	50:DT:112:ARG:HG3	2.31	0.60
50:DT:28:VAL:O	50:DT:88:ILE:HD11	2.00	0.60
51:DU:66:ASN:OD1	51:DU:76:TYR:N	2.34	0.60
54:DX:61:GLY:H	54:DX:70:LEU:HD21	1.66	0.60
55:DY:28:LYS:CA	55:DY:39:VAL:H	2.15	0.60
55:DY:68:HIS:HB3	55:DY:71:LYS:CE	2.32	0.60
55:DY:81:LYS:HG2	55:DY:97:ARG:H	1.65	0.60
56:DZ:77:ASP:OD1	56:DZ:79:ARG:HD2	2.00	0.60
1:AA:824:C:H4'	8:AH:1:MET:N	2.16	0.60
3:AC:53:ALA:HB2	3:AC:115:LEU:CD2	2.29	0.60
5:AE:122:GLU:O	5:AE:123:LEU:HD23	2.02	0.60
7:AG:95:ARG:O	7:AG:96:GLN:C	2.40	0.60
7:AG:93:PRO:HA	7:AG:96:GLN:NE2	2.16	0.60
23:AW:50:G:H2'	23:AW:51:U:O4'	2.01	0.60
29:B3:52:HIS:CD2	36:BB:83:G:H4'	2.37	0.60
29:B3:56:VAL:O	29:B3:57:GLU:HB2	2.02	0.60
31:B5:25:LEU:CD1	53:BW:19:LEU:HB3	2.32	0.60
34:B8:43:GLN:O	34:B8:44:LYS:HD2	2.01	0.60
35:BA:2103:C:C2'	35:BA:2104:G:H5''	2.32	0.60
35:BA:540:C:H2'	35:BA:541:C:C5	2.36	0.60
35:BA:580:C:H2'	35:BA:581:C:C6	2.37	0.60
37:BC:58:VAL:HG21	37:BC:166:ASP:N	2.15	0.60
38:BD:209:ALA:C	38:BD:210:GLY:O	2.37	0.60
39:BE:52:LEU:O	39:BE:74:PRO:HA	2.01	0.60
39:BE:77:ILE:HG21	39:BE:79:ARG:HE	1.66	0.60
40:BF:22:ALA:O	40:BF:26:ALA:HB2	2.01	0.60
42:BH:122:THR:O	42:BH:133:VAL:HG13	2.00	0.60
42:BH:12:PRO:O	42:BH:13:LYS:HB2	2.01	0.60
44:BN:15:LEU:HD12	44:BN:136:GLU:HB2	1.83	0.60
45:BO:79:PHE:CE2	45:BO:101:PRO:HB2	2.32	0.60
45:BO:31:LYS:HD2	45:BO:32:TYR:HE1	1.66	0.60
46:BP:107:LYS:C	46:BP:109:GLY:H	2.04	0.60
46:BP:33:ARG:O	46:BP:34:GLY:C	2.39	0.60
47:BQ:52:VAL:CG1	47:BQ:53:ALA:H	2.05	0.60
48:BR:84:ALA:HB3	48:BR:85:PRO:HD3	1.84	0.60
45:BO:77:ILE:HD11	50:BT:72:VAL:CG1	2.32	0.60
53:BW:87:PRO:HA	53:BW:93:ALA:CB	2.32	0.60
54:BX:92:LEU:O	54:BX:93:GLU:HB3	2.02	0.60
55:BY:89:PHE:C	55:BY:90:LEU:HD23	2.23	0.60
1:CA:1168:A:H2'	1:CA:1169:A:C8	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1281:U:H5'	1:CA:1282:C:H5	1.66	0.60
1:CA:1391:U:H2'	1:CA:1392:G:H8	1.66	0.60
1:CA:321:A:H4'	1:CA:1436:U:O4'	2.02	0.60
1:CA:445:G:H2'	1:CA:446:G:H8	1.67	0.60
1:CA:634:C:O2'	1:CA:635:G:H5'	2.02	0.60
1:CA:707:C:H4'	11:CK:20:TYR:HD1	1.66	0.60
1:CA:731:G:OP1	1:CA:766:A:H1'	2.02	0.60
2:CB:158:LEU:HD12	2:CB:158:LEU:N	2.17	0.60
2:CB:80:ILE:HG21	2:CB:208:ILE:HG23	1.82	0.60
4:CD:126:ILE:HG22	4:CD:127:THR:H	1.65	0.60
4:CD:79:PHE:O	4:CD:82:ALA:HB3	2.01	0.60
5:CE:6:PHE:HB3	5:CE:35:GLY:O	2.02	0.60
8:CH:34:GLU:O	8:CH:38:ILE:HG13	2.02	0.60
14:CN:33:VAL:HA	14:CN:39:LEU:O	2.02	0.60
15:CO:18:PHE:CZ	15:CO:21:ASP:HB2	2.37	0.60
19:CS:6:LYS:HD2	19:CS:7:LYS:HD2	1.84	0.60
23:CW:19:G:H3'	23:CW:20:G:C5'	2.31	0.60
25:CY:133:ARG:NH1	35:DA:1942:C:O4'	2.34	0.60
28:D2:29:LYS:O	28:D2:33:MET:N	2.31	0.60
35:DA:1327:C:H2'	35:DA:1328:G:O4'	2.02	0.60
35:DA:141:A:C8	35:DA:1408:C:O2'	2.54	0.60
35:DA:2092:U:C5	35:DA:2226:C:OP2	2.53	0.60
35:DA:84:A:N3	35:DA:85:G:H1'	2.16	0.60
38:DD:223:GLY:O	38:DD:225:ALA:N	2.29	0.60
35:DA:2579:C:O3'	39:DE:131:ALA:HB2	2.02	0.60
39:DE:176:ILE:HG22	39:DE:179:GLU:H	1.65	0.60
41:DG:162:THR:O	41:DG:164:GLU:N	2.35	0.60
42:DH:158:HIS:NE2	42:DH:170:ARG:HA	2.17	0.60
44:DN:16:ILE:O	44:DN:54:VAL:HA	2.01	0.60
44:DN:82:LEU:HD12	44:DN:83:LYS:N	2.17	0.60
46:DP:46:LYS:HB3	46:DP:52:GLU:HG2	1.83	0.60
50:DT:100:TYR:O	50:DT:102:ILE:N	2.35	0.60
56:DZ:11:GLU:OE2	56:DZ:12:GLY:N	2.35	0.60
1:AA:101:A:H2'	1:AA:102:G:H8	1.67	0.59
1:AA:989:C:N4	1:AA:1216:G:H1	1.99	0.59
1:AA:190:U:H2'	1:AA:191:G:H8	1.66	0.59
2:AB:16:HIS:HA	2:AB:210:SER:HB2	1.84	0.59
4:AD:106:TYR:CE1	4:AD:113:SER:HA	2.37	0.59
4:AD:163:GLU:C	4:AD:165:MET:H	2.04	0.59
4:AD:173:TRP:C	4:AD:186:LEU:HD12	2.21	0.59
8:AH:22:GLU:O	8:AH:63:LEU:HD23	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:AX:14:U:H2'	24:AX:15:G:C8	2.37	0.59
27:B1:71:TYR:HA	27:B1:74:VAL:HG23	1.82	0.59
28:B2:32:LEU:HG	28:B2:33:MET:N	2.17	0.59
31:B5:20:ARG:HA	31:B5:23:HIS:HD2	1.67	0.59
35:BA:1678:G:N2	35:BA:1989:G:H22	1.99	0.59
35:BA:2575:C:H5'	39:BE:144:ARG:HG2	1.84	0.59
38:BD:175:LEU:HD23	38:BD:176:ARG:N	2.16	0.59
38:BD:231:HIS:CE1	38:BD:232:PRO:HD2	2.36	0.59
39:BE:29:GLY:HA3	39:BE:180:ASN:ND2	2.16	0.59
44:BN:62:VAL:HG22	44:BN:66:LYS:CG	2.31	0.59
46:BP:130:PHE:N	46:BP:130:PHE:HD2	1.99	0.59
46:BP:21:ARG:HH11	46:BP:21:ARG:HG3	1.67	0.59
47:BQ:22:LYS:NZ	47:BQ:22:LYS:HA	2.17	0.59
52:BV:37:VAL:HG12	52:BV:38:LEU:N	2.16	0.59
54:BX:36:LYS:HD3	54:BX:38:GLU:HB2	1.83	0.59
55:BY:19:LYS:HD2	55:BY:20:TYR:CE1	2.37	0.59
1:CA:1118:C:OP1	9:CI:9:ARG:HD3	2.01	0.59
1:CA:1246:C:H2'	1:CA:1247:U:C6	2.36	0.59
1:CA:620:C:H2'	1:CA:621:A:O4'	2.00	0.59
1:CA:93:G:O2'	1:CA:96:U:H5'	2.02	0.59
5:CE:69:VAL:O	5:CE:71:LEU:N	2.35	0.59
7:CG:74:GLU:H	7:CG:91:VAL:HG23	1.66	0.59
13:CM:90:LEU:C	13:CM:92:HIS:N	2.50	0.59
20:CT:73:HIS:HB3	20:CT:74:LYS:CD	2.28	0.59
31:D5:32:PRO:O	31:D5:33:CYS:HB2	2.03	0.59
35:DA:116:C:O2'	35:DA:117:G:H5'	2.02	0.59
35:DA:1190:G:O5'	46:DP:35:HIS:HA	2.02	0.59
35:DA:1385:G:H4'	35:DA:1386:C:OP1	2.01	0.59
35:DA:1486:A:H61	35:DA:1504:C:H42	1.50	0.59
35:DA:2514:U:H2'	35:DA:2515:C:H6	1.67	0.59
35:DA:2586:C:O2'	35:DA:2587:A:H5'	2.02	0.59
35:DA:2810:A:H2'	39:DE:61:ARG:HH21	1.66	0.59
39:DE:52:LEU:O	39:DE:74:PRO:HA	2.02	0.59
39:DE:9:VAL:HG22	39:DE:25:VAL:HB	1.84	0.59
40:DF:161:GLU:O	40:DF:164:ARG:HB2	2.02	0.59
40:DF:200:GLU:O	40:DF:204:ASN:HB2	2.02	0.59
40:DF:202:PHE:HD1	40:DF:202:PHE:C	2.05	0.59
41:DG:132:ASN:CB	41:DG:159:VAL:HG22	2.26	0.59
44:DN:78:TYR:H	44:DN:79:PRO:HD2	1.67	0.59
46:DP:30:THR:CG2	46:DP:31:ALA:H	2.08	0.59
47:DQ:27:VAL:HG23	56:DZ:81:ARG:HH22	1.64	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:89:PHE:C	55:DY:90:LEU:HD23	2.23	0.59
56:DZ:118:GLN:HA	56:DZ:118:GLN:HE21	1.67	0.59
1:AA:1088:G:H2'	1:AA:1089:G:C8	2.35	0.59
1:AA:176:C:H2'	1:AA:177:C:H6	1.67	0.59
1:AA:262:A:H2'	1:AA:263:A:C8	2.38	0.59
1:AA:503:C:H2'	1:AA:504:C:H6	1.67	0.59
1:AA:694:A:H5''	11:AK:53:SER:CB	2.32	0.59
7:AG:100:ALA:C	7:AG:104:LEU:HD23	2.23	0.59
8:AH:2:LEU:O	8:AH:3:THR:HG23	2.01	0.59
25:AY:150:SER:HB2	25:AY:153:GLU:H	1.67	0.59
25:AY:15:GLN:O	25:AY:18:LEU:HB3	2.02	0.59
25:AY:3:LEU:H	25:AY:3:LEU:CD1	2.13	0.59
25:AY:65:THR:HA	25:AY:103:ILE:HG23	1.84	0.59
29:B3:56:VAL:O	29:B3:57:GLU:CB	2.50	0.59
33:B7:16:HIS:HD1	33:B7:21:ARG:HH22	1.50	0.59
35:BA:1158:C:HO2'	35:BA:1159:U:H6	1.50	0.59
35:BA:1215:G:H2'	35:BA:1216:G:H8	1.67	0.59
35:BA:1438:U:H2'	35:BA:1439:A:H8	1.66	0.59
35:BA:1484:G:H3'	35:BA:1485:G:H5''	1.82	0.59
35:BA:2092:U:C5	35:BA:2226:C:OP2	2.54	0.59
35:BA:2752:C:H5'	35:BA:2753:A:OP2	2.02	0.59
35:BA:2843:G:H1	35:BA:2874:C:H42	1.50	0.59
35:BA:1789:A:OP1	38:BD:222:ARG:HG3	2.02	0.59
38:BD:68:LYS:HB2	38:BD:70:TRP:CH2	2.37	0.59
39:BE:23:VAL:HA	39:BE:184:VAL:O	2.02	0.59
41:BG:173:LEU:N	41:BG:173:LEU:HD22	2.17	0.59
46:BP:126:VAL:HA	46:BP:145:PRO:CB	2.31	0.59
35:BA:1190:G:C5'	46:BP:35:HIS:HA	2.33	0.59
47:BQ:140:ALA:CB	56:BZ:99:TYR:HB2	2.33	0.59
47:BQ:22:LYS:HA	47:BQ:22:LYS:HZ3	1.66	0.59
51:BU:58:ARG:HA	51:BU:61:TRP:CE3	2.37	0.59
52:BV:72:VAL:HG12	52:BV:73:SER:N	2.13	0.59
54:BX:49:VAL:CG1	54:BX:50:LYS:H	2.12	0.59
56:BZ:68:PRO:HG2	56:BZ:91:LEU:O	2.01	0.59
1:CA:431:A:H2'	1:CA:432:A:H8	1.67	0.59
1:CA:688:G:O2'	1:CA:689:C:H5'	2.02	0.59
1:CA:922:G:H2'	1:CA:923:A:C8	2.36	0.59
3:CC:109:PRO:HA	3:CC:115:LEU:HD13	1.83	0.59
3:CC:111:LEU:CD2	3:CC:146:ALA:HB2	2.32	0.59
6:CF:91:VAL:HG11	18:CR:72:ARG:NH1	2.16	0.59
7:CG:143:ARG:HH11	7:CG:143:ARG:CB	2.15	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CJ:50:ILE:N	10:CJ:50:ILE:HD13	2.00	0.59
35:DA:214:G:O2'	35:DA:215:G:O4'	2.19	0.59
36:DB:114:C:H4'	49:DS:46:VAL:HG13	1.84	0.59
39:DE:35:GLN:HE22	39:DE:37:ARG:NH2	2.00	0.59
39:DE:3:GLY:O	39:DE:4:ILE:HG22	2.02	0.59
41:DG:13:GLU:O	41:DG:14:GLU:HB3	2.02	0.59
41:DG:135:LEU:HD23	41:DG:155:MET:CE	2.31	0.59
42:DH:70:THR:CG2	42:DH:74:ASN:HD21	2.13	0.59
46:DP:106:LEU:HD11	46:DP:112:LEU:HB2	1.84	0.59
46:DP:56:SER:O	46:DP:57:THR:HB	2.02	0.59
46:DP:92:GLU:HG3	46:DP:93:GLY:H	1.66	0.59
47:DQ:22:LYS:HZ3	47:DQ:22:LYS:HA	1.67	0.59
53:DW:20:VAL:CG2	53:DW:47:VAL:HG21	2.31	0.59
54:DX:72:LYS:CE	54:DX:74:PRO:HB3	2.24	0.59
56:DZ:114:GLY:O	56:DZ:115:GLY:O	2.21	0.59
1:AA:1014:A:H2'	1:AA:1015:A:C8	2.36	0.59
1:AA:1281:U:H5'	1:AA:1282:C:H5	1.67	0.59
1:AA:123:C:OP1	1:AA:312:C:H5'	2.02	0.59
1:AA:473:G:H2'	1:AA:474:G:H8	1.67	0.59
1:AA:922:G:H2'	1:AA:923:A:C8	2.37	0.59
3:AC:153:VAL:HA	3:AC:197:GLY:O	2.02	0.59
4:AD:118:ARG:O	4:AD:121:VAL:HB	2.02	0.59
5:AE:80:ILE:CG1	5:AE:91:LEU:HB2	2.32	0.59
17:AQ:9:VAL:HG12	17:AQ:10:VAL:N	2.18	0.59
20:AT:51:GLU:O	20:AT:55:ILE:HG12	2.01	0.59
21:AU:2:GLY:N	21:AU:5:ASP:HB2	2.18	0.59
31:B5:8:LYS:O	31:B5:9:LYS:HD2	2.03	0.59
35:BA:1446:C:N4	35:BA:1465:G:H1	1.98	0.59
35:BA:1509(A):A:H2'	35:BA:1509(B):A:C8	2.36	0.59
35:BA:173:G:H2'	35:BA:174:C:C6	2.38	0.59
35:BA:1803:A:H4'	38:BD:259:THR:HG23	1.84	0.59
35:BA:2103:C:H2'	35:BA:2104:G:H5''	1.83	0.59
35:BA:2732:G:H3'	35:BA:2733:A:C5'	2.33	0.59
35:BA:816:C:H2'	35:BA:817:C:H6	1.65	0.59
35:BA:979:G:H3'	35:BA:980:A:C5'	2.31	0.59
39:BE:81:ILE:O	39:BE:81:ILE:HG22	2.02	0.59
40:BF:50:SER:HB2	40:BF:94:PRO:HD3	1.84	0.59
44:BN:132:ALA:O	44:BN:133:GLN:CB	2.50	0.59
44:BN:44:PRO:C	44:BN:46:VAL:H	2.04	0.59
46:BP:127:ALA:HB3	46:BP:130:PHE:HE2	1.61	0.59
46:BP:16:ARG:NE	46:BP:18:ARG:HB2	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BR:37:THR:CG2	48:BR:40:LYS:HE2	2.32	0.59
50:BT:100:TYR:O	50:BT:102:ILE:N	2.35	0.59
51:BU:45:TYR:O	51:BU:46:ALA:C	2.41	0.59
51:BU:83:LEU:CG	51:BU:88:ILE:HG12	2.32	0.59
52:BV:40:LEU:O	52:BV:41:GLY:C	2.41	0.59
54:BX:89:ILE:HD12	54:BX:89:ILE:N	2.17	0.59
56:BZ:105:VAL:H	56:BZ:141:VAL:HG11	1.67	0.59
56:BZ:163:LEU:HD23	56:BZ:163:LEU:N	2.16	0.59
1:CA:1045:C:H2'	1:CA:1046:A:O4'	2.02	0.59
1:CA:1056:U:H5'	3:CC:163:ALA:HB2	1.84	0.59
1:CA:1498:U:H1'	1:CA:1499:A:OP2	2.01	0.59
1:CA:445:G:H2'	1:CA:446:G:C8	2.37	0.59
1:CA:943:U:H2'	1:CA:944:G:C8	2.35	0.59
3:CC:92:ALA:N	3:CC:99:VAL:HG21	2.16	0.59
10:CJ:5:ARG:HG3	10:CJ:73:ASP:OD1	2.01	0.59
11:CK:87:THR:HA	11:CK:91:ARG:HG3	1.85	0.59
12:CL:85:ILE:HD11	12:CL:98:TYR:HB2	1.83	0.59
1:CA:375:U:H4'	16:CP:17:TYR:CE2	2.37	0.59
25:CY:108:GLU:O	25:CY:112:LYS:N	2.28	0.59
27:D1:46:LEU:HD22	27:D1:48:LYS:HE2	1.84	0.59
35:DA:1232:G:H2'	35:DA:1233:C:C6	2.36	0.59
35:DA:1366:A:H2'	35:DA:1367:A:C8	2.37	0.59
35:DA:141:A:H8	35:DA:1408:C:HO2'	1.49	0.59
35:DA:2590:A:O2'	35:DA:2591:C:H5'	2.02	0.59
35:DA:30:G:O2'	35:DA:31:C:H5'	2.02	0.59
35:DA:42:G:H2'	35:DA:43:A:C8	2.36	0.59
35:DA:796:C:O2'	35:DA:797:C:H5'	2.01	0.59
35:DA:869:G:H1'	47:DQ:8:LYS:HZ2	1.66	0.59
39:DE:55:ASN:HD21	39:DE:75:VAL:HG22	1.66	0.59
40:DF:199:TRP:O	40:DF:203:GLN:HG2	2.02	0.59
40:DF:88:VAL:CG2	40:DF:89:VAL:N	2.65	0.59
42:DH:46:GLU:O	42:DH:47:GLU:HB2	2.01	0.59
43:DI:76:THR:HB	43:DI:139:GLN:O	2.02	0.59
44:DN:34:LEU:HD23	44:DN:120:LEU:HD23	1.84	0.59
45:DO:13:ASN:HD22	45:DO:97:ARG:CG	2.15	0.59
34:D8:30:ARG:HE	46:DP:62:LEU:HB2	1.66	0.59
48:DR:13:HIS:O	48:DR:14:SER:O	2.20	0.59
48:DR:28:LEU:HA	48:DR:34:ILE:CG1	2.32	0.59
51:DU:30:LYS:HE3	51:DU:30:LYS:HA	1.85	0.59
55:DY:86:ARG:HG2	55:DY:87:LYS:N	2.18	0.59
1:AA:1035:A:H2'	1:AA:1036:G:H8	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1478:C:H2'	1:AA:1479:C:C5	2.37	0.59
1:AA:280:C:C2	17:AQ:38:ARG:HG3	2.37	0.59
1:AA:861:G:H2'	1:AA:862:C:H6	1.67	0.59
1:AA:943:U:H2'	1:AA:944:G:C8	2.32	0.59
4:AD:18:LYS:HE3	4:AD:31:CYS:SG	2.43	0.59
6:AF:30:LEU:HD23	6:AF:75:LEU:HD21	1.84	0.59
13:AM:82:MET:HB3	13:AM:93:ARG:HH11	1.67	0.59
1:AA:375:U:H4'	16:AP:17:TYR:CE2	2.38	0.59
16:AP:71:ARG:HA	16:AP:74:LEU:HD12	1.83	0.59
20:AT:63:ILE:O	20:AT:65:LYS:N	2.35	0.59
27:B1:41:ARG:NH2	35:BA:205:G:H1	2.00	0.59
34:B8:36:LYS:O	34:B8:37:SER:O	2.19	0.59
35:BA:1177:A:H5''	35:BA:1178:C:O5'	2.03	0.59
35:BA:1794:U:O4'	35:BA:1900:A:C2	2.56	0.59
35:BA:2266:A:H4'	35:BA:2267:A:N3	2.17	0.59
35:BA:2468:G:HO2'	35:BA:2476:A:H8	1.51	0.59
35:BA:2061:G:H5''	35:BA:2503:A:C2	2.37	0.59
35:BA:422:A:H2'	35:BA:423:A:C8	2.36	0.59
35:BA:1792:G:P	38:BD:206:LEU:HB2	2.42	0.59
38:BD:94:LEU:H	38:BD:94:LEU:CD1	2.14	0.59
39:BE:36:ARG:NH2	39:BE:88:GLY:CA	2.66	0.59
42:BH:87:LEU:C	42:BH:88:LEU:HD22	2.22	0.59
47:BQ:121:ALA:O	47:BQ:125:LEU:HD12	2.02	0.59
49:BS:84:GLN:HA	49:BS:105:ALA:O	2.02	0.59
49:BS:72:ALA:O	49:BS:76:LYS:HG2	2.01	0.59
39:BE:14:ILE:HB	50:BT:14:TYR:CE2	2.37	0.59
53:BW:20:VAL:CG2	53:BW:47:VAL:HG21	2.31	0.59
54:BX:39:ILE:O	54:BX:42:ALA:HB3	2.02	0.59
54:BX:40:LYS:C	54:BX:42:ALA:H	2.04	0.59
54:BX:81:VAL:HG12	54:BX:82:GLN:O	2.02	0.59
56:BZ:148:ASP:O	56:BZ:149:SER:HB3	2.01	0.59
1:CA:451:A:H1'	1:CA:452:A:C8	2.37	0.59
1:CA:59:A:H3'	1:CA:331:G:H22	1.68	0.59
1:CA:67:C:H2'	1:CA:68:G:H8	1.65	0.59
1:CA:909:A:C2	1:CA:910:C:H1'	2.38	0.59
6:CF:21:LEU:HA	6:CF:24:GLU:HG2	1.85	0.59
11:CK:61:ALA:CB	11:CK:90:GLY:HA3	2.31	0.59
15:CO:23:GLY:O	15:CO:24:SER:HB3	2.02	0.59
15:CO:24:SER:O	15:CO:28:GLN:HG3	2.02	0.59
18:CR:56:THR:OG1	18:CR:57:GLY:N	2.35	0.59
35:DA:1453:U:H5'	48:DR:63:ARG:NE	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:268:C:H42	35:DA:424:G:H1	1.48	0.59
35:DA:2795:G:C2	35:DA:2799:C:H5'	2.37	0.59
37:DC:45:ALA:H	37:DC:174:PRO:CB	2.14	0.59
38:DD:31:LYS:HZ1	38:DD:31:LYS:HA	1.68	0.59
50:DT:94:ALA:HB1	50:DT:99:LEU:HD23	1.84	0.59
51:DU:83:LEU:CG	51:DU:88:ILE:HG12	2.32	0.59
54:DX:88:LYS:HD2	54:DX:88:LYS:N	2.16	0.59
56:DZ:25:PRO:HB2	56:DZ:85:HIS:ND1	2.17	0.59
56:DZ:99:TYR:HE2	56:DZ:125:LEU:HD12	1.67	0.59
1:AA:1416:G:H2'	1:AA:1417:G:H8	1.67	0.59
1:AA:445:G:H2'	1:AA:446:G:C8	2.37	0.59
1:AA:447:G:N2	1:AA:488:C:H42	1.99	0.59
1:AA:528:C:H2'	1:AA:529:G:H8	1.67	0.59
1:AA:552:U:O3'	12:AL:87:GLY:HA3	2.03	0.59
1:AA:923:A:H2'	1:AA:924:C:C6	2.38	0.59
1:AA:927:G:OP2	1:AA:1503:A:C4	2.56	0.59
2:AB:163:PHE:HA	2:AB:185:ILE:O	2.02	0.59
2:AB:24:TRP:CG	2:AB:25:ASN:N	2.69	0.59
3:AC:130:VAL:O	3:AC:134:ILE:HG13	2.01	0.59
4:AD:180:GLY:C	4:AD:181:MET:HG2	2.22	0.59
7:AG:74:GLU:H	7:AG:91:VAL:HG23	1.67	0.59
19:AS:10:PHE:HE2	19:AS:70:LYS:HZ1	1.51	0.59
23:AW:60:A:H2'	23:AW:61:U:O4'	2.02	0.59
23:AW:68:C:H2'	23:AW:69:C:O4'	2.02	0.59
25:AY:61:PRO:CG	25:AY:67:VAL:HG13	2.31	0.59
31:B5:2:ALA:N	35:BA:747:U:N3	2.49	0.59
31:B5:30:LEU:HD11	53:BW:38:TYR:HB2	1.84	0.59
32:B6:14:THR:O	32:B6:49:HIS:HA	2.03	0.59
35:BA:1198:U:H2'	35:BA:1199:U:C6	2.37	0.59
35:BA:1217:C:H2'	35:BA:1218:C:O4'	2.03	0.59
35:BA:1362:C:O2'	35:BA:1363:C:H5'	2.02	0.59
35:BA:1418:G:N1	35:BA:1579:A:H5'	2.16	0.59
35:BA:142:A:H5'	35:BA:142(A):C:OP2	2.03	0.59
35:BA:2279:G:N2	35:BA:2280:G:H1'	2.17	0.59
35:BA:2320:A:H8	35:BA:2321:G:O6	1.86	0.59
35:BA:2624:G:O2'	35:BA:2625:G:H5'	2.03	0.59
35:BA:2744:G:N7	35:BA:2755:C:O2	2.35	0.59
35:BA:700:G:H2'	35:BA:701:G:C8	2.38	0.59
35:BA:700:G:H2'	35:BA:701:G:H8	1.68	0.59
39:BE:176:ILE:HG22	39:BE:179:GLU:H	1.67	0.59
43:BI:31:LEU:HB3	43:BI:32:PRO:HD3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:99:LEU:O	50:BT:99:LEU:CD1	2.50	0.59
53:BW:17:VAL:O	53:BW:19:LEU:N	2.34	0.59
28:B2:33:MET:HG2	54:BX:10:ALA:CB	2.32	0.59
54:BX:51:VAL:HG13	54:BX:80:ILE:N	2.18	0.59
55:BY:16:ALA:C	55:BY:21:LYS:HD2	2.23	0.59
56:BZ:19:ARG:HB3	56:BZ:19:ARG:NH1	2.16	0.59
1:CA:190:U:H2'	1:CA:191:G:H8	1.68	0.59
1:CA:383:A:H8	1:CA:383:A:O5'	1.84	0.59
1:CA:834:C:H2'	1:CA:835:U:H6	1.67	0.59
1:CA:865:A:H2'	1:CA:866:C:O4'	2.01	0.59
1:CA:975:A:H4'	1:CA:976:G:C5'	2.26	0.59
2:CB:51:LEU:HB3	2:CB:55:PHE:HE2	1.67	0.59
4:CD:11:LEU:O	4:CD:13:ARG:N	2.36	0.59
5:CE:146:ALA:C	5:CE:148:VAL:H	2.06	0.59
6:CF:82:ARG:HB2	6:CF:85:VAL:CG2	2.32	0.59
13:CM:9:ILE:CG2	13:CM:11:ARG:HG3	2.32	0.59
17:CQ:68:ARG:N	17:CQ:70:ARG:HH12	1.99	0.59
20:CT:14:LYS:HA	20:CT:17:ARG:HH21	1.66	0.59
28:D2:57:ILE:HG13	28:D2:57:ILE:O	2.02	0.59
35:DA:1330:C:O2'	35:DA:1331:A:H5'	2.03	0.59
35:DA:1344:G:H4'	35:DA:1384:A:C5	2.38	0.59
35:DA:1505:C:H2'	35:DA:1506:C:O4'	2.03	0.59
35:DA:1754:C:OP1	50:DT:96:ARG:NH1	2.34	0.59
35:DA:1844:C:C2'	35:DA:1845:G:H5'	2.31	0.59
35:DA:1992:G:C6	35:DA:1997:G:N1	2.71	0.59
35:DA:2744:G:O2'	35:DA:2745:C:H5'	2.03	0.59
35:DA:2787:C:C2	39:DE:61:ARG:HD3	2.37	0.59
35:DA:407:G:H2'	35:DA:408:G:H8	1.67	0.59
38:DD:206:LEU:HD23	38:DD:211:ARG:HH11	1.68	0.59
38:DD:81:ALA:N	38:DD:94:LEU:HD11	2.16	0.59
40:DF:114:VAL:HG11	40:DF:202:PHE:HE2	1.67	0.59
41:DG:139:LEU:O	41:DG:144:ILE:HG21	2.03	0.59
41:DG:56:ALA:HB1	41:DG:153:ARG:HD2	1.83	0.59
36:DB:42:C:O4'	41:DG:69:ALA:HB2	2.03	0.59
42:DH:144:VAL:O	42:DH:144:VAL:HG12	2.01	0.59
44:DN:23:LEU:HB3	44:DN:60:ILE:CG2	2.32	0.59
44:DN:9:VAL:HG12	44:DN:10:GLU:N	2.15	0.59
45:DO:77:ILE:HD11	50:DT:72:VAL:CG1	2.33	0.59
46:DP:13:ASN:HD22	46:DP:13:ASN:H	1.51	0.59
46:DP:144:GLU:N	46:DP:145:PRO:HD3	2.17	0.59
47:DQ:29:PHE:HB2	47:DQ:65:PHE:CE2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1278:A:H5''	48:DR:36:THR:HG22	1.84	0.59
54:DX:82:GLN:CG	54:DX:83:VAL:N	2.65	0.59
1:AA:1030(D):A:H2'	1:AA:1031:G:H5'	1.85	0.59
1:AA:1442(A):G:N2	50:BT:119:LYS:N	2.50	0.59
1:AA:192:U:H2'	1:AA:193:C:H6	1.68	0.59
1:AA:370:C:O2'	1:AA:371:G:H5'	2.03	0.59
1:AA:403:C:O2'	1:AA:404:U:H5'	2.03	0.59
4:AD:127:THR:HG23	4:AD:147:ALA:HB3	1.85	0.59
4:AD:3:ARG:O	4:AD:5:ILE:N	2.36	0.59
4:AD:65:ARG:HB2	4:AD:75:PHE:CE2	2.37	0.59
5:AE:131:ILE:HD13	5:AE:131:ILE:N	2.17	0.59
10:AJ:32:ALA:N	10:AJ:78:ASN:ND2	2.51	0.59
10:AJ:30:SER:OG	10:AJ:81:THR:HG22	2.03	0.59
12:AL:90:VAL:HG11	12:AL:93:LEU:HG	1.84	0.59
23:AW:23:G:C2'	23:AW:24:C:H5''	2.32	0.59
27:B1:73:LEU:CA	27:B1:76:ARG:NH1	2.63	0.59
29:B3:4:LEU:HD23	29:B3:5:LYS:N	2.17	0.59
35:BA:1264:G:H3'	35:BA:1265:A:H5''	1.84	0.59
35:BA:1778:U:C5	35:BA:1784:A:C4	2.91	0.59
35:BA:2248:C:C2'	35:BA:2249:U:H5'	2.32	0.59
35:BA:2569:G:O2'	35:BA:2570:G:H5'	2.02	0.59
35:BA:2669:G:H2'	35:BA:2670:A:H8	1.68	0.59
35:BA:270:A:O2'	35:BA:271:A:H5'	2.03	0.59
35:BA:2762:G:C2'	35:BA:2763:G:H5''	2.32	0.59
35:BA:2766:G:N3	35:BA:2766:G:H2'	2.15	0.59
35:BA:481:G:H1'	35:BA:506:G:H21	1.68	0.59
35:BA:601:C:O2	35:BA:605:C:H4'	2.01	0.59
38:BD:176:ARG:HG2	38:BD:176:ARG:HH11	1.66	0.59
38:BD:215:LEU:O	38:BD:217:ARG:N	2.35	0.59
39:BE:92:THR:HB	39:BE:94:GLU:OE1	2.03	0.59
41:BG:57:ALA:HA	41:BG:60:LEU:HB3	1.84	0.59
42:BH:68:THR:HA	42:BH:71:LEU:HD22	1.84	0.59
44:BN:56:ASN:HA	44:BN:125:GLY:H	1.68	0.59
47:BQ:29:PHE:CD1	47:BQ:29:PHE:N	2.71	0.59
50:BT:128:GLU:O	50:BT:130:ALA:N	2.35	0.59
51:BU:31:SER:HB3	51:BU:34:LYS:HB2	1.84	0.59
51:BU:33:ARG:O	51:BU:34:LYS:C	2.41	0.59
52:BV:2:PHE:CB	52:BV:42:GLY:HA2	2.29	0.59
52:BV:79:VAL:HG12	52:BV:80:GLN:N	2.17	0.59
1:CA:1438:G:N2	1:CA:1464:G:H1'	2.18	0.59
1:CA:390:C:H4'	16:CP:28:ARG:HH21	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:591:U:H2'	1:CA:592:G:H8	1.65	0.59
1:CA:642:A:N3	8:CH:113:SER:OG	2.26	0.59
1:CA:949:A:H61	1:CA:1232:U:H3	1.51	0.59
1:CA:473:G:OP1	16:CP:81:ARG:HB2	2.02	0.59
27:D1:76:ARG:CB	27:D1:78:LYS:HZ3	2.15	0.59
27:D1:85:LEU:N	27:D1:85:LEU:HD23	2.17	0.59
31:D5:2:ALA:N	35:DA:2014:A:N3	2.50	0.59
35:DA:1180:C:H2'	35:DA:1181:C:H5'	1.83	0.59
35:DA:1289:C:H2'	35:DA:1290:C:C6	2.37	0.59
35:DA:2121:G:H1	35:DA:2177:C:H42	1.49	0.59
35:DA:319:C:O2'	35:DA:320:A:H5'	2.03	0.59
35:DA:644:A:C2	35:DA:2369:A:H1'	2.38	0.59
35:DA:669:G:H2'	35:DA:669:G:N3	2.17	0.59
37:DC:44:HIS:HA	37:DC:174:PRO:CB	2.32	0.59
38:DD:10:THR:HG23	38:DD:13:ARG:HB3	1.84	0.59
38:DD:134:ARG:HB2	38:DD:135:PHE:HD1	1.66	0.59
38:DD:142:VAL:HG22	38:DD:143:HIS:N	2.18	0.59
38:DD:160:GLY:H	38:DD:196:VAL:HB	1.67	0.59
39:DE:179:GLU:HB3	39:DE:181:LEU:CD2	2.26	0.59
41:DG:83:ARG:O	41:DG:85:GLY:N	2.36	0.59
44:DN:26:LEU:HG	44:DN:30:ILE:HD11	1.84	0.59
46:DP:100:LEU:HD22	46:DP:100:LEU:N	2.17	0.59
35:DA:598:G:H5'	46:DP:15:ARG:CD	2.30	0.59
47:DQ:83:MET:O	47:DQ:83:MET:CG	2.51	0.59
50:DT:28:VAL:O	50:DT:28:VAL:HG12	2.03	0.59
50:DT:78:LEU:O	50:DT:79:HIS:ND1	2.36	0.59
51:DU:106:PHE:CA	51:DU:109:LEU:HD12	2.29	0.59
53:DW:35:ILE:HG22	53:DW:36:LEU:HD23	1.85	0.59
53:DW:74:ALA:O	53:DW:75:TYR:HB3	2.03	0.59
54:DX:82:GLN:HG3	54:DX:83:VAL:N	2.17	0.59
1:AA:1152:A:H2'	1:AA:1153:C:H6	1.67	0.59
1:AA:1205:U:O2'	3:AC:195:VAL:HG23	2.01	0.59
1:AA:1264:C:H2'	1:AA:1265:G:C8	2.37	0.59
1:AA:1292:U:O2'	1:AA:1293:G:H5'	2.01	0.59
1:AA:512:U:H2'	1:AA:513:C:C6	2.38	0.59
1:AA:764:C:H2'	1:AA:765:G:C8	2.37	0.59
1:AA:786:G:H1	1:AA:796:C:H42	1.49	0.59
2:AB:36:ARG:NH1	2:AB:37:ASN:HB2	2.17	0.59
2:AB:69:LEU:HB2	2:AB:159:PRO:HG2	1.83	0.59
5:AE:37:ARG:O	5:AE:38:GLN:HG2	2.02	0.59
9:AI:37:PHE:CE1	9:AI:74:ILE:HG12	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:109:VAL:HG22	18:AR:86:VAL:HA	1.85	0.59
1:AA:363:A:C2	12:AL:31:PRO:HG2	2.38	0.59
27:B1:27:GLU:H	27:B1:34:THR:HB	1.67	0.59
35:BA:1503:U:H2'	35:BA:1504:C:C5	2.37	0.59
35:BA:1613:G:H2'	35:BA:1617:C:H42	1.66	0.59
35:BA:1804:C:O2'	35:BA:1805:U:H5'	2.02	0.59
35:BA:2039:C:H2'	35:BA:2040:C:H6	1.68	0.59
35:BA:2261:C:H1'	35:BA:2388:A:N3	2.17	0.59
35:BA:2772:C:H2'	35:BA:2773:C:H6	1.67	0.59
31:B5:29:THR:HG21	35:BA:2814:C:O2'	2.03	0.59
35:BA:686:G:N2	35:BA:788:A:N6	2.49	0.59
39:BE:110:GLY:O	48:BR:2:ARG:CZ	2.50	0.59
39:BE:36:ARG:HH22	39:BE:88:GLY:H	1.47	0.59
39:BE:52:LEU:CB	39:BE:76:ARG:HB2	2.28	0.59
41:BG:121:ASN:HB2	41:BG:181:ARG:HH21	1.66	0.59
45:BO:122:LEU:N	45:BO:122:LEU:HD12	2.17	0.59
50:BT:23:ARG:HG2	50:BT:120:ARG:HH12	1.66	0.59
50:BT:14:TYR:N	50:BT:14:TYR:HD1	2.00	0.59
51:BU:59:ARG:O	51:BU:61:TRP:N	2.35	0.59
52:BV:22:VAL:O	52:BV:23:GLU:HB2	2.03	0.59
54:BX:12:VAL:CG1	54:BX:27:THR:HG23	2.33	0.59
55:BY:2:ARG:C	55:BY:4:LYS:H	2.05	0.59
56:BZ:55:HIS:CE1	56:BZ:133:ILE:HG21	2.38	0.59
56:BZ:58:VAL:CG2	56:BZ:68:PRO:HA	2.30	0.59
2:CB:69:LEU:HB3	2:CB:162:ILE:HG22	1.83	0.59
4:CD:62:GLN:HB3	4:CD:66:ARG:NH1	2.18	0.59
6:CF:82:ARG:HB3	6:CF:82:ARG:NH1	2.18	0.59
8:CH:127:LEU:HD13	8:CH:127:LEU:O	2.01	0.59
8:CH:27:PRO:HA	8:CH:58:TYR:HA	1.83	0.59
8:CH:45:ILE:HA	8:CH:64:LYS:CB	2.32	0.59
1:CA:1349:A:OP1	9:CI:120:ARG:HB3	2.02	0.59
10:CJ:79:ARG:HA	10:CJ:82:ILE:HG12	1.84	0.59
12:CL:115:LYS:O	12:CL:117:ARG:HG3	2.03	0.59
15:CO:82:ILE:HD11	15:CO:87:ILE:O	2.03	0.59
31:D5:22:HIS:N	31:D5:22:HIS:ND1	2.51	0.59
32:D6:42:TRP:HA	32:D6:42:TRP:HE3	1.68	0.59
33:D7:7:PRO:CB	35:DA:1309:G:H4'	2.33	0.59
34:D8:32:LEU:HD11	34:D8:41:ILE:CG2	2.32	0.59
35:DA:1264:G:H3'	35:DA:1265:A:H5''	1.84	0.59
35:DA:1827:C:O2'	35:DA:1828:G:H5'	2.02	0.59
35:DA:2327:A:H2'	35:DA:2328:A:C8	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2525:G:H2'	35:DA:2526:G:H8	1.67	0.59
35:DA:2529:G:OP2	35:DA:2530:A:H5''	2.03	0.59
35:DA:2704:C:C2'	35:DA:2705:A:H8	2.15	0.59
35:DA:2774:C:H2'	35:DA:2775:A:C8	2.38	0.59
35:DA:27:G:N2	35:DA:512:G:C2'	2.64	0.59
35:DA:805:G:H22	35:DA:828:U:H5''	1.68	0.59
38:DD:186:HIS:CD2	38:DD:188:GLU:H	2.12	0.59
40:DF:50:SER:HB2	40:DF:94:PRO:HD3	1.84	0.59
41:DG:150:ASP:O	41:DG:151:ALA:HB2	2.02	0.59
44:DN:132:ALA:O	44:DN:133:GLN:CB	2.51	0.59
44:DN:1:MET:C	44:DN:2:LYS:HD2	2.23	0.59
44:DN:99:LEU:O	44:DN:103:VAL:HG23	2.02	0.59
45:DO:37:ASP:H	45:DO:62:VAL:H	1.50	0.59
45:DO:43:VAL:O	45:DO:45:GLU:N	2.36	0.59
48:DR:20:LEU:HD12	48:DR:21:TYR:N	2.17	0.59
50:DT:92:GLY:C	50:DT:94:ALA:N	2.56	0.59
51:DU:110:VAL:O	51:DU:114:LYS:N	2.32	0.59
52:DV:14:VAL:HG12	52:DV:15:GLU:N	2.18	0.59
1:AA:1483:A:C2'	1:AA:1484:C:H5'	2.33	0.59
1:AA:555:C:H2'	1:AA:556:C:C6	2.38	0.59
1:AA:967:C:H2'	1:AA:968:A:N7	2.18	0.59
2:AB:167:PRO:HG2	2:AB:192:SER:OG	2.03	0.59
3:AC:60:ALA:HB3	3:AC:63:ASN:HD21	1.67	0.59
4:AD:108:LEU:HD11	4:AD:174:LEU:HD22	1.84	0.59
4:AD:61:LYS:HE3	4:AD:207:TYR:OH	2.02	0.59
5:AE:12:LEU:H	5:AE:12:LEU:HD13	1.67	0.59
6:AF:71:ARG:O	6:AF:73:ASN:N	2.36	0.59
11:AK:99:GLN:C	11:AK:101:SER:H	2.06	0.59
12:AL:32:PHE:HB3	12:AL:84:LEU:HD21	1.84	0.59
16:AP:2:VAL:HG23	16:AP:22:THR:O	2.03	0.59
17:AQ:44:ALA:HB1	17:AQ:73:VAL:HG22	1.84	0.59
25:AY:140:LEU:HD11	25:AY:157:ALA:CB	2.33	0.59
27:B1:58:ILE:HD11	27:B1:87:PRO:HB3	1.84	0.59
35:BA:1505:C:H2'	35:BA:1506:C:O4'	2.02	0.59
35:BA:1688:U:H1'	35:BA:1701:A:C6	2.38	0.59
35:BA:1799:G:H4'	35:BA:1800:C:O5'	2.02	0.59
35:BA:1935:G:H1'	35:BA:1964:G:N2	2.17	0.59
35:BA:2208:A:H1'	35:BA:2219:G:C2	2.38	0.59
35:BA:2525:G:H2'	35:BA:2526:G:H8	1.68	0.59
35:BA:58:G:H1	35:BA:69:C:H42	1.50	0.59
35:BA:807:U:H2'	35:BA:808:G:C8	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1792:G:OP2	38:BD:206:LEU:HD12	2.03	0.59
41:BG:61:ALA:HA	41:BG:64:THR:HG22	1.85	0.59
48:BR:37:THR:OG1	48:BR:40:LYS:HG3	2.03	0.59
49:BS:87:PHE:O	49:BS:88:ASP:HB2	2.03	0.59
50:BT:129:ARG:CZ	50:BT:131:ALA:HB3	2.32	0.59
55:BY:83:THR:HG22	55:BY:84:ARG:N	2.18	0.59
56:BZ:103:ARG:HE	56:BZ:103:ARG:HA	1.67	0.59
1:CA:1320:C:H5'	19:CS:70:LYS:HE3	1.84	0.59
1:CA:659:U:H2'	1:CA:660:G:H8	1.67	0.59
5:CE:11:ILE:HD12	5:CE:31:LEU:HD22	1.85	0.59
8:CH:45:ILE:HB	8:CH:62:TYR:O	2.03	0.59
9:CI:118:LYS:HB3	9:CI:118:LYS:HZ3	1.66	0.59
12:CL:22:SER:C	12:CL:24:VAL:H	2.06	0.59
17:CQ:44:ALA:HB1	17:CQ:73:VAL:HG22	1.85	0.59
35:DA:1703:G:H2'	35:DA:1704:G:H8	1.67	0.59
35:DA:257:A:H2'	35:DA:258:G:O4'	2.03	0.59
37:DC:211:SER:HA	37:DC:220:PRO:HA	1.85	0.59
35:DA:1812:A:H1'	38:DD:46:GLN:HE22	1.67	0.59
39:DE:200:GLU:OE2	39:DE:200:GLU:N	2.35	0.59
41:DG:40:ASN:HD22	41:DG:91:ARG:CG	2.16	0.59
41:DG:76:SER:HB3	41:DG:83:ARG:HA	1.85	0.59
42:DH:13:LYS:O	42:DH:15:VAL:N	2.35	0.59
43:DI:102:SER:HA	43:DI:107:VAL:O	2.03	0.59
43:DI:68:LEU:CD2	43:DI:136:VAL:HG11	2.32	0.59
44:DN:40:PRO:HG3	51:DU:68:ALA:HB2	1.84	0.59
47:DQ:81:VAL:HG23	47:DQ:82:ARG:HH11	1.67	0.59
49:DS:65:VAL:O	49:DS:69:VAL:HG12	2.03	0.59
50:DT:128:GLU:O	50:DT:130:ALA:N	2.35	0.59
50:DT:14:TYR:CD1	50:DT:14:TYR:N	2.67	0.59
51:DU:98:LEU:O	51:DU:101:ARG:N	2.36	0.59
51:DU:47:TYR:HA	51:DU:50:ARG:NH1	2.17	0.59
51:DU:91:ASP:O	51:DU:95:LEU:HB2	2.03	0.59
52:DV:24:LYS:HA	52:DV:94:LEU:HG	1.84	0.59
35:DA:565:C:O3'	52:DV:81:TYR:CE1	2.56	0.59
56:DZ:148:ASP:O	56:DZ:149:SER:HB3	2.03	0.59
56:DZ:149:SER:HB2	56:DZ:172:ALA:O	2.03	0.59
1:AA:1109:C:H2'	1:AA:1110:A:O4'	2.03	0.59
1:AA:1294:G:H2'	1:AA:1295:G:C8	2.38	0.59
1:AA:600:C:H5'	8:AH:129:VAL:O	2.03	0.59
1:AA:811:C:O2'	1:AA:901:A:N1	2.36	0.59
1:AA:974:A:OP1	14:AN:31:ARG:HD3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:107:THR:HG23	2:AB:110:GLN:OE1	2.03	0.59
2:AB:25:ASN:OD1	2:AB:27:LYS:HB2	2.03	0.59
6:AF:10:LEU:HD13	6:AF:61:LEU:HD11	1.83	0.59
7:AG:22:LEU:O	7:AG:25:ALA:HB3	2.03	0.59
10:AJ:50:ILE:CD1	10:AJ:50:ILE:H	1.91	0.59
25:AY:63:PRO:CB	25:AY:64:ARG:HH22	2.15	0.59
27:B1:44:PRO:HA	35:BA:2231:C:OP1	2.02	0.59
28:B2:29:LYS:HA	28:B2:32:LEU:CD2	2.30	0.59
32:B6:15:GLU:O	32:B6:16:CYS:SG	2.59	0.59
34:B8:32:LEU:HD11	34:B8:41:ILE:CG2	2.33	0.59
35:BA:1824:G:O2'	35:BA:1825:A:H5'	2.02	0.59
35:BA:2024:G:H2'	35:BA:2025:C:H6	1.66	0.59
35:BA:2774:C:H2'	35:BA:2775:A:H8	1.66	0.59
35:BA:514:A:H1'	35:BA:581:C:O2'	2.03	0.59
35:BA:64:A:H2'	35:BA:65:C:H6	1.66	0.59
29:B3:14:GLY:O	35:BA:969:U:H4'	2.02	0.59
38:BD:142:VAL:HG23	38:BD:193:VAL:CA	2.31	0.59
39:BE:55:ASN:HD21	39:BE:75:VAL:HG22	1.66	0.59
41:BG:83:ARG:O	41:BG:85:GLY:N	2.36	0.59
43:BI:37:VAL:CG1	43:BI:38:LEU:N	2.65	0.59
44:BN:46:VAL:HG22	44:BN:47:ALA:N	2.17	0.59
45:BO:24:VAL:HG21	45:BO:32:TYR:O	2.03	0.59
45:BO:63:VAL:HG22	45:BO:83:ALA:O	2.03	0.59
45:BO:93:PRO:C	45:BO:95:GLY:H	2.04	0.59
50:BT:27:THR:OG1	50:BT:28:VAL:N	2.36	0.59
54:BX:52:VAL:O	54:BX:53:LYS:CB	2.51	0.59
55:BY:86:ARG:HG2	55:BY:87:LYS:N	2.18	0.59
56:BZ:166:SER:HB2	56:BZ:168:GLU:N	2.18	0.59
1:CA:1292:U:H2'	1:CA:1293:G:H8	1.68	0.59
1:CA:135:C:H2'	1:CA:136:C:H5'	1.83	0.59
1:CA:370:C:O2'	1:CA:371:G:H5'	2.03	0.59
1:CA:974:A:OP1	14:CN:31:ARG:HD3	2.03	0.59
2:CB:121:LEU:O	2:CB:121:LEU:HD23	2.03	0.59
2:CB:73:THR:HG22	2:CB:93:VAL:O	2.03	0.59
3:CC:88:ARG:HG2	3:CC:101:LEU:HB2	1.84	0.59
4:CD:154:ASN:CB	4:CD:159:ARG:HH21	2.15	0.59
4:CD:196:LEU:HD12	4:CD:196:LEU:N	2.18	0.59
7:CG:36:LYS:HB2	7:CG:36:LYS:NZ	2.17	0.59
18:CR:86:VAL:O	18:CR:87:ARG:HB3	2.01	0.59
19:CS:16:LEU:CD1	19:CS:16:LEU:H	2.15	0.59
25:CY:171:LYS:O	25:CY:172:ALA:C	2.40	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:D8:60:LEU:HA	34:D8:63:PRO:CG	2.33	0.59
35:DA:1127:A:C2'	35:DA:1128:A:H5''	2.33	0.59
35:DA:1555:G:H2'	35:DA:1556:C:C6	2.38	0.59
35:DA:1789:A:OP1	38:DD:222:ARG:HG3	2.03	0.59
35:DA:493:G:H2'	35:DA:494:G:H5''	1.85	0.59
35:DA:674:G:H1'	40:DF:74:ARG:HG3	1.85	0.59
35:DA:700:G:H2'	35:DA:701:G:C8	2.37	0.59
37:DC:184:LYS:C	37:DC:186:ALA:H	2.04	0.59
38:DD:132:PRO:HD3	38:DD:190:TYR:CZ	2.37	0.59
44:DN:15:LEU:HD12	44:DN:136:GLU:HB2	1.84	0.59
44:DN:46:VAL:O	44:DN:47:ALA:HB2	2.03	0.59
45:DO:46:ALA:H	45:DO:54:GLU:HG2	1.66	0.59
47:DQ:22:LYS:HA	47:DQ:22:LYS:NZ	2.18	0.59
50:DT:28:VAL:N	50:DT:88:ILE:HD13	2.17	0.59
50:DT:80:SER:O	50:DT:82:LEU:N	2.36	0.59
55:DY:43:ASN:O	55:DY:44:ILE:O	2.21	0.59
1:AA:1074:G:H4'	2:AB:103:THR:HG22	1.85	0.59
4:AD:155:LEU:O	4:AD:158:ILE:HG22	2.03	0.59
4:AD:60:GLU:HG2	4:AD:202:LEU:HB2	1.84	0.59
16:AP:19:ILE:HB	16:AP:37:GLY:CA	2.32	0.59
18:AR:36:ASN:HD22	18:AR:39:VAL:CB	2.16	0.59
18:AR:44:LEU:HA	18:AR:49:LYS:O	2.03	0.59
18:AR:87:ARG:CZ	18:AR:87:ARG:HB3	2.33	0.59
23:AW:57:C:H2'	23:AW:58:A:H8	1.67	0.59
28:B2:12:GLU:C	28:B2:14:ARG:NE	2.57	0.59
35:BA:1374:G:H2'	35:BA:1375:C:C6	2.38	0.59
35:BA:1937:A:N7	35:BA:1939:U:H2'	2.18	0.59
35:BA:2892:A:C5	35:BA:2893:G:H1'	2.38	0.59
35:BA:742:G:O2'	35:BA:743:G:H5'	2.01	0.59
38:BD:95:LEU:HD12	38:BD:103:ARG:O	2.03	0.59
38:BD:270:ILE:HD12	38:BD:270:ILE:O	2.03	0.59
38:BD:32:SER:OG	38:BD:33:LEU:N	2.34	0.59
35:BA:2579:C:O3'	39:BE:131:ALA:HB2	2.03	0.59
39:BE:75:VAL:C	39:BE:77:ILE:H	2.06	0.59
42:BH:13:LYS:O	42:BH:15:VAL:N	2.35	0.59
43:BI:14:ASP:HB2	43:BI:17:GLN:OE1	2.03	0.59
44:BN:53:VAL:HG13	44:BN:121:LYS:O	2.02	0.59
46:BP:71:VAL:HG22	46:BP:72:PRO:CD	2.33	0.59
48:BR:49:ASP:O	48:BR:50:HIS:C	2.41	0.59
51:BU:47:TYR:HA	51:BU:50:ARG:CZ	2.32	0.59
1:CA:1152:A:H2'	1:CA:1153:C:H6	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1228:C:H4'	13:CM:116:THR:O	2.02	0.59
3:CC:87:LEU:HB3	3:CC:101:LEU:HD11	1.85	0.59
7:CG:149:ARG:HB3	11:CK:59:TYR:CE2	2.37	0.59
8:CH:9:MET:O	8:CH:13:ILE:HG12	2.02	0.59
8:CH:91:ARG:HG2	8:CH:91:ARG:HH11	1.68	0.59
10:CJ:30:SER:OG	10:CJ:81:THR:HG22	2.02	0.59
1:CA:972:C:H2'	10:CJ:55:LYS:HD3	1.84	0.59
12:CL:32:PHE:HB3	12:CL:84:LEU:CD2	2.33	0.59
12:CL:89:ARG:C	12:CL:89:ARG:HH11	2.07	0.59
29:D3:51:ALA:O	29:D3:53:LEU:N	2.36	0.59
34:D8:60:LEU:HA	34:D8:63:PRO:HG2	1.84	0.59
35:DA:1438:U:H2'	35:DA:1439:A:H8	1.66	0.59
35:DA:1630:G:H2'	35:DA:1631:C:H6	1.68	0.59
35:DA:1720:U:C2'	35:DA:1721:G:H5'	2.32	0.59
35:DA:2469:A:H3'	35:DA:2470:G:O4'	2.02	0.59
35:DA:2669:G:H2'	35:DA:2670:A:H8	1.68	0.59
35:DA:2673:G:O2'	35:DA:2674:G:H5'	2.03	0.59
35:DA:2762:G:C2'	35:DA:2763:G:H5''	2.30	0.59
35:DA:2872:G:C2	35:DA:2873:A:N6	2.70	0.59
35:DA:377:C:H2'	35:DA:378:C:H6	1.64	0.59
35:DA:37:C:H2'	35:DA:38:A:C8	2.38	0.59
35:DA:437:G:O2'	35:DA:438:G:H5'	2.03	0.59
35:DA:700:G:H2'	35:DA:701:G:H8	1.67	0.59
38:DD:80:ALA:HB2	38:DD:96:HIS:ND1	2.17	0.59
39:DE:103:ASP:OD2	39:DE:202:LYS:HE2	2.03	0.59
35:DA:2682:U:C2	39:DE:22:PRO:HB3	2.38	0.59
39:DE:77:ILE:HG21	39:DE:79:ARG:HE	1.67	0.59
43:DI:109:ILE:CD1	43:DI:111:PRO:HD3	2.33	0.59
43:DI:95:LYS:O	43:DI:99:GLU:HB2	2.03	0.59
44:DN:26:LEU:HD11	44:DN:30:ILE:HD11	1.85	0.59
44:DN:55:VAL:HG12	44:DN:56:ASN:N	2.18	0.59
46:DP:108:LYS:O	46:DP:110:TYR:N	2.32	0.59
51:DU:91:ASP:OD2	51:DU:96:ALA:CA	2.51	0.59
1:AA:9:G:H2'	1:AA:10:A:H8	1.68	0.58
1:AA:1241:G:H2'	1:AA:1242:C:C5	2.38	0.58
1:AA:949:A:H2'	1:AA:950:U:O4'	2.03	0.58
5:AE:15:ARG:O	5:AE:15:ARG:HG2	2.03	0.58
8:AH:11:THR:HA	8:AH:14:ARG:CZ	2.33	0.58
9:AI:95:LYS:NZ	9:AI:96:LEU:HD13	2.18	0.58
10:AJ:79:ARG:HA	10:AJ:82:ILE:HG12	1.84	0.58
14:AN:33:VAL:HA	14:AN:39:LEU:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:AS:41:VAL:CB	19:AS:44:MET:HB2	2.33	0.58
25:AY:3:LEU:HD23	25:AY:7:TYR:OH	2.03	0.58
31:B5:2:ALA:N	35:BA:2014:A:N3	2.51	0.58
35:BA:1180:C:H2'	35:BA:1181:C:H5'	1.85	0.58
35:BA:407:G:H2'	35:BA:408:G:H8	1.68	0.58
35:BA:13:A:C2	35:BA:526:A:C5	2.90	0.58
35:BA:680:G:H2'	35:BA:681:G:C8	2.37	0.58
38:BD:142:VAL:HA	38:BD:194:GLY:H	1.68	0.58
39:BE:131:ALA:HB1	39:BE:134:ILE:HD11	1.85	0.58
39:BE:169:ASN:OD1	39:BE:201:THR:HG21	2.03	0.58
41:BG:64:THR:HG23	41:BG:65:GLY:N	2.17	0.58
45:BO:68:GLU:HB3	45:BO:78:ARG:HB2	1.85	0.58
35:BA:1245:G:C3'	46:BP:16:ARG:HH22	2.15	0.58
50:BT:28:VAL:O	50:BT:28:VAL:HG12	2.02	0.58
50:BT:89:VAL:HG11	50:BT:91:ARG:HE	1.67	0.58
51:BU:6:THR:HG21	51:BU:10:ARG:NH2	2.18	0.58
51:BU:34:LYS:HA	51:BU:34:LYS:CE	2.32	0.58
54:BX:30:VAL:HG23	54:BX:76:ARG:HA	1.85	0.58
56:BZ:33:LEU:HD12	56:BZ:34:ASN:H	1.68	0.58
56:BZ:63:ASP:C	56:BZ:65:GLN:H	2.07	0.58
1:CA:599:C:H2'	1:CA:600:C:H6	1.68	0.58
1:CA:745:C:H2'	1:CA:746:A:H8	1.68	0.58
2:CB:31:TYR:HD2	2:CB:31:TYR:N	2.02	0.58
4:CD:163:GLU:C	4:CD:165:MET:H	2.06	0.58
5:CE:37:ARG:O	5:CE:38:GLN:HG2	2.02	0.58
9:CI:37:PHE:CE1	9:CI:74:ILE:HG12	2.38	0.58
1:CA:972:C:H5'	10:CJ:57:LYS:NZ	2.17	0.58
11:CK:22:HIS:C	11:CK:28:THR:HG23	2.22	0.58
25:CY:76:LEU:HD23	25:CY:76:LEU:C	2.24	0.58
25:CY:41:LEU:HD22	25:CY:83:ILE:HD13	1.84	0.58
27:D1:13:ILE:CG2	27:D1:14:VAL:H	2.12	0.58
27:D1:89:GLU:OE2	27:D1:89:GLU:N	2.35	0.58
28:D2:17:SER:O	28:D2:18:PRO:C	2.39	0.58
29:D3:35:ARG:HG2	29:D3:37:LEU:HD21	1.85	0.58
35:DA:1227:G:OP2	51:DU:16:LYS:HE3	2.03	0.58
35:DA:1309:G:O2'	35:DA:1310:G:H5'	2.02	0.58
35:DA:176:G:C2'	35:DA:177:G:H5'	2.33	0.58
35:DA:2051:A:H5'	35:DA:2578:G:O4'	2.02	0.58
35:DA:575:A:O2'	35:DA:576:U:H5'	2.03	0.58
35:DA:692:C:H2'	35:DA:693:C:C6	2.38	0.58
38:DD:132:PRO:HG3	38:DD:190:TYR:CE1	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:231:HIS:CD2	38:DD:249:PRO:HG3	2.38	0.58
39:DE:29:GLY:HA3	39:DE:180:ASN:ND2	2.16	0.58
40:DF:3:GLU:HA	40:DF:24:LEU:HB3	1.85	0.58
41:DG:109:VAL:O	41:DG:113:ARG:N	2.36	0.58
44:DN:42:TRP:HD1	51:DU:63:VAL:HG11	1.67	0.58
50:DT:28:VAL:CG1	50:DT:46:GLU:HA	2.32	0.58
53:DW:58:ALA:HB1	53:DW:64:MET:SD	2.43	0.58
54:DX:65:ARG:NH2	54:DX:66:LEU:H	2.00	0.58
56:DZ:109:ALA:O	56:DZ:111:VAL:HG12	2.03	0.58
56:DZ:27:VAL:HG13	56:DZ:27:VAL:O	2.02	0.58
1:AA:1152:A:H5''	10:AJ:13:HIS:HD2	1.66	0.58
1:AA:1368:G:O2'	1:AA:1369:C:H5'	2.03	0.58
1:AA:909:A:C2	1:AA:910:C:H1'	2.37	0.58
2:AB:71:VAL:O	2:AB:164:VAL:HA	2.02	0.58
1:AA:1075:C:OP1	2:AB:179:LYS:HE2	2.03	0.58
2:AB:63:MET:HB3	2:AB:225:ALA:HB1	1.85	0.58
4:AD:119:GLN:HG3	4:AD:123:HIS:CD2	2.38	0.58
4:AD:18:LYS:NZ	4:AD:31:CYS:HB3	2.18	0.58
5:AE:57:LYS:HE2	5:AE:61:TYR:HE2	1.67	0.58
8:AH:110:ALA:CB	8:AH:121:ASP:HB3	2.26	0.58
9:AI:28:VAL:HG13	9:AI:64:THR:CA	2.29	0.58
11:AK:29:ILE:HD12	11:AK:29:ILE:C	2.23	0.58
7:AG:153:HIS:CE1	11:AK:57:THR:HG23	2.39	0.58
15:AO:8:LYS:O	15:AO:12:ILE:HG13	2.03	0.58
16:AP:19:ILE:HG22	16:AP:36:ILE:CG1	2.33	0.58
1:AA:390:C:H4'	16:AP:28:ARG:HH21	1.68	0.58
17:AQ:86:GLU:C	17:AQ:88:TYR:N	2.56	0.58
18:AR:86:VAL:O	18:AR:87:ARG:HB3	2.02	0.58
19:AS:22:LEU:HD22	19:AS:27:GLU:H	1.69	0.58
19:AS:63:THR:HG22	19:AS:66:MET:CG	2.25	0.58
25:AY:29:ARG:NE	25:AY:32:ARG:NH2	2.44	0.58
25:AY:39:LEU:HB2	25:AY:53:ASN:CB	2.20	0.58
34:B8:30:ARG:HE	46:BP:62:LEU:HB2	1.68	0.58
34:B8:48:PHE:H	34:B8:48:PHE:HD1	1.50	0.58
35:BA:2078:C:H2'	35:BA:2079:U:H6	1.68	0.58
35:BA:2408:U:H2'	35:BA:2409:G:H8	1.68	0.58
35:BA:2455:G:H2'	35:BA:2456:C:C6	2.38	0.58
35:BA:271(Q):G:H2'	35:BA:271(R):G:C8	2.38	0.58
35:BA:2723:C:H2'	35:BA:2724:C:H5'	1.84	0.58
35:BA:590:A:H2'	35:BA:591:C:C6	2.38	0.58
35:BA:8:A:H2'	35:BA:9:U:C6	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BC:196:LEU:C	37:BC:198:ALA:H	2.07	0.58
39:BE:30:PRO:C	39:BE:32:PRO:HD3	2.23	0.58
35:BA:2636:U:H4'	39:BE:80:GLU:OE1	2.01	0.58
35:BA:607:U:OP1	40:BF:102:PRO:HA	2.02	0.58
40:BF:200:GLU:O	40:BF:204:ASN:HB2	2.02	0.58
40:BF:71:GLY:O	40:BF:72:ARG:C	2.41	0.58
41:BG:137:GLU:HB3	41:BG:140:ILE:CG2	2.32	0.58
42:BH:140:LYS:O	42:BH:144:VAL:HG23	2.03	0.58
44:BN:23:LEU:HB3	44:BN:60:ILE:CG2	2.33	0.58
44:BN:65:LYS:CE	44:BN:65:LYS:HA	2.27	0.58
44:BN:77:GLY:O	44:BN:78:TYR:HB3	2.03	0.58
46:BP:122:PRO:CG	46:BP:141:ALA:HB3	2.33	0.58
47:BQ:25:ASP:HA	56:BZ:78:LYS:HZ2	1.65	0.58
48:BR:74:LYS:O	48:BR:77:ARG:N	2.36	0.58
54:BX:34:ALA:O	54:BX:36:LYS:HG3	2.03	0.58
54:BX:61:GLY:H	54:BX:70:LEU:HD21	1.68	0.58
55:BY:28:LYS:HE3	55:BY:28:LYS:O	2.02	0.58
56:BZ:120:ILE:HG21	56:BZ:170:THR:O	2.02	0.58
1:CA:690:G:H2'	1:CA:691:G:O4'	2.03	0.58
1:CA:959:A:H2'	1:CA:960:U:H4'	1.85	0.58
4:CD:3:ARG:O	4:CD:5:ILE:N	2.37	0.58
10:CJ:48:THR:HG22	10:CJ:49:VAL:N	2.18	0.58
10:CJ:4:ILE:HB	10:CJ:74:ILE:CG1	2.32	0.58
11:CK:19:ALA:HA	11:CK:32:ILE:HA	1.85	0.58
16:CP:5:ARG:C	16:CP:6:LEU:HD12	2.24	0.58
16:CP:71:ARG:HH11	16:CP:71:ARG:HG3	1.68	0.58
18:CR:74:ARG:HA	18:CR:79:LEU:HB2	1.84	0.58
19:CS:40:ILE:HD13	19:CS:62:ILE:HD13	1.85	0.58
23:CW:71:G:O2'	23:CW:72:C:H5'	2.02	0.58
25:CY:13:HIS:H	25:CY:13:HIS:HD1	1.50	0.58
26:D0:43:THR:HB	26:D0:57:PHE:CE1	2.38	0.58
27:D1:87:PRO:O	27:D1:90:ILE:HG12	2.03	0.58
29:D3:31:LEU:CD2	29:D3:32:GLN:H	2.15	0.58
32:D6:27:LYS:HD2	35:DA:2285:C:OP2	2.03	0.58
35:DA:1809:A:H2'	35:DA:1810:A:C8	2.38	0.58
35:DA:1862:G:O2'	35:DA:1863:G:H5'	2.02	0.58
35:DA:2626:C:H2'	35:DA:2627:G:C8	2.37	0.58
35:DA:2780:G:OP1	44:DN:118:LYS:HE2	2.04	0.58
35:DA:2808:U:O2'	35:DA:2809:A:H5'	2.03	0.58
35:DA:361:G:C2'	35:DA:362:U:H5''	2.33	0.58
35:DA:360:G:H2'	35:DA:361:G:H8	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:575:A:H2'	35:DA:576:U:H5'	1.83	0.58
35:DA:94(A):G:C3'	35:DA:95:G:H5''	2.33	0.58
35:DA:2632:A:O2'	39:DE:61:ARG:NH2	2.36	0.58
43:DI:133:HIS:CB	43:DI:134:PRO:CD	2.81	0.58
47:DQ:35:VAL:CG1	47:DQ:130:LYS:HB3	2.33	0.58
45:DO:76:ALA:HB3	50:DT:75:ILE:CB	2.32	0.58
51:DU:92:ARG:HB3	52:DV:11:GLN:CD	2.21	0.58
53:DW:29:LEU:O	53:DW:33:ARG:HG3	2.02	0.58
54:DX:36:LYS:HZ1	54:DX:39:ILE:HA	1.67	0.58
54:DX:40:LYS:C	54:DX:42:ALA:H	2.07	0.58
54:DX:68:ARG:HG3	54:DX:69:TYR:CD1	2.37	0.58
56:DZ:121:HIS:C	56:DZ:123:ASP:H	2.05	0.58
1:AA:1228:C:H4'	13:AM:116:THR:O	2.03	0.58
1:AA:1363(A):A:H4'	1:AA:1364:U:C5'	2.24	0.58
1:AA:338:A:H2'	1:AA:339:C:H6	1.68	0.58
1:AA:533:A:H1'	1:AA:534:U:OP1	2.03	0.58
1:AA:745:C:H2'	1:AA:746:A:H8	1.68	0.58
2:AB:69:LEU:CB	2:AB:162:ILE:HG22	2.34	0.58
2:AB:98:LEU:HB2	2:AB:101:MET:CG	2.34	0.58
3:AC:109:PRO:HA	3:AC:115:LEU:HD13	1.85	0.58
11:AK:80:VAL:HG23	11:AK:80:VAL:O	2.03	0.58
13:AM:68:GLY:O	13:AM:69:GLU:HB2	2.02	0.58
23:AW:59:A:O3'	23:AW:61:U:H5	1.85	0.58
25:AY:29:ARG:NH2	25:AY:110:ARG:HH21	2.00	0.58
29:B3:15:TYR:HB3	29:B3:19:GLN:NE2	2.17	0.58
35:BA:1050:A:O2'	35:BA:2752:C:H1'	2.02	0.58
35:BA:122:G:H1	35:BA:129:C:H42	1.51	0.58
35:BA:812:C:C2	35:BA:1250:G:N1	2.71	0.58
35:BA:1623:G:H2'	35:BA:1624:G:H8	1.69	0.58
35:BA:1658:C:OP1	39:BE:132:HIS:ND1	2.35	0.58
35:BA:1887:C:C3'	35:BA:1888:G:H5''	2.33	0.58
27:B1:41:ARG:HH12	35:BA:189:G:H3'	1.68	0.58
35:BA:2024:G:O2'	35:BA:2025:C:H5'	2.02	0.58
35:BA:271(L):U:H4'	35:BA:271(M):G:C5	2.38	0.58
38:BD:130:ALA:HB2	38:BD:192:THR:CB	2.33	0.58
38:BD:201:HIS:O	38:BD:203:ASN:N	2.36	0.58
38:BD:79:VAL:HG11	38:BD:112:GLN:O	2.04	0.58
35:BA:2729:G:C1'	39:BE:187:ALA:HB2	2.21	0.58
41:BG:173:LEU:O	41:BG:176:LEU:HB2	2.03	0.58
42:BH:144:VAL:O	42:BH:144:VAL:HG12	2.02	0.58
44:BN:1:MET:HG2	44:BN:2:LYS:N	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:81:GLN:NE2	46:BP:106:LEU:HA	2.18	0.58
49:BS:23:ARG:HG2	49:BS:24:LEU:H	1.67	0.58
53:BW:1:MET:HE3	53:BW:1:MET:HA	1.85	0.58
1:CA:1368:G:O2'	1:CA:1369:C:H5'	2.03	0.58
1:CA:1439:C:O2	1:CA:1439:C:H2'	2.03	0.58
1:CA:259:G:H2'	1:CA:260:G:C8	2.38	0.58
1:CA:393:A:O2'	1:CA:394:G:H5'	2.03	0.58
1:CA:452:A:O2'	1:CA:453:A:H8	1.75	0.58
1:CA:627:G:H2'	1:CA:628:G:C8	2.38	0.58
2:CB:98:LEU:HB2	2:CB:101:MET:CG	2.32	0.58
4:CD:173:TRP:C	4:CD:186:LEU:HD12	2.23	0.58
1:CA:1298:C:C4	7:CG:114:ARG:HD2	2.38	0.58
7:CG:152:ALA:O	7:CG:154:TYR:N	2.36	0.58
8:CH:103:VAL:HG21	8:CH:109:ILE:C	2.22	0.58
8:CH:110:ALA:CB	8:CH:121:ASP:HB3	2.27	0.58
9:CI:14:VAL:O	9:CI:65:VAL:HG23	2.02	0.58
9:CI:47:LEU:HB3	9:CI:50:LEU:HD12	1.84	0.58
16:CP:72:ARG:C	16:CP:74:LEU:H	2.06	0.58
25:CY:15:GLN:HA	25:CY:168:PHE:HZ	1.66	0.58
27:D1:73:LEU:O	27:D1:76:ARG:HG2	2.02	0.58
28:D2:12:GLU:O	28:D2:14:ARG:CZ	2.51	0.58
29:D3:3:ARG:HA	29:D3:38:GLU:HA	1.85	0.58
31:D5:57:VAL:HG23	31:D5:58:LEU:N	2.19	0.58
34:D8:25:MET:HG3	46:DP:62:LEU:HD21	1.85	0.58
34:D8:42:ARG:O	34:D8:44:LYS:N	2.36	0.58
35:DA:1217:C:H2'	35:DA:1218:C:O4'	2.03	0.58
35:DA:2078:C:H2'	35:DA:2079:U:H6	1.69	0.58
35:DA:2320:A:H8	35:DA:2321:G:O6	1.86	0.58
35:DA:2627:G:N3	35:DA:2781:A:H2	2.02	0.58
35:DA:2808:U:H2'	35:DA:2809:A:C5'	2.34	0.58
35:DA:742:G:O2'	35:DA:743:G:H5'	2.03	0.58
35:DA:827:U:H2'	35:DA:2068:U:N3	2.17	0.58
35:DA:851:U:H2'	35:DA:852:G:C8	2.35	0.58
37:DC:68:LEU:HB3	37:DC:70:LYS:HG2	1.85	0.58
40:DF:202:PHE:CD1	40:DF:202:PHE:C	2.76	0.58
40:DF:65:TRP:CH2	40:DF:75:HIS:CD2	2.87	0.58
41:DG:46:ALA:C	41:DG:51:ARG:HG3	2.23	0.58
41:DG:95:ARG:O	41:DG:99:MET:N	2.30	0.58
43:DI:115:ALA:O	43:DI:128:LEU:HD23	2.03	0.58
45:DO:46:ALA:N	45:DO:54:GLU:HG2	2.18	0.58
46:DP:39:LYS:C	46:DP:41:ARG:H	2.05	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:62:LEU:C	46:DP:62:LEU:HD22	2.23	0.58
46:DP:64:LYS:O	46:DP:65:ARG:C	2.42	0.58
47:DQ:124:LYS:HE2	47:DQ:124:LYS:HA	1.85	0.58
49:DS:61:ASN:ND2	49:DS:62:LYS:HE3	2.14	0.58
51:DU:6:THR:HG21	51:DU:10:ARG:NH2	2.18	0.58
53:DW:9:TYR:HD2	53:DW:102:HIS:HE2	1.49	0.58
1:AA:1313:U:OP1	19:AS:6:LYS:HG3	2.03	0.58
1:AA:1343:G:H1'	9:AI:121:ARG:HH12	1.69	0.58
1:AA:1391:U:H2'	1:AA:1392:G:H8	1.65	0.58
1:AA:1397:C:N4	24:AX:22:U:C5	2.71	0.58
3:AC:91:LEU:O	3:AC:94:LEU:HG	2.03	0.58
4:AD:79:PHE:O	4:AD:82:ALA:HB3	2.03	0.58
6:AF:41:GLU:H	6:AF:62:TRP:HE3	1.52	0.58
7:AG:79:ARG:NH2	23:AW:34:U:H4'	2.18	0.58
10:AJ:49:VAL:HG13	14:AN:41:ARG:HB2	1.84	0.58
12:AL:25:PRO:O	12:AL:27:LEU:HD22	2.03	0.58
12:AL:5:PRO:HG2	12:AL:10:LEU:HD21	1.83	0.58
12:AL:6:THR:HG22	12:AL:9:GLN:CG	2.31	0.58
13:AM:117:VAL:HG12	13:AM:118:ALA:N	2.17	0.58
27:B1:13:ILE:HG23	27:B1:14:VAL:N	2.18	0.58
31:B5:44:THR:CG2	31:B5:45:VAL:H	2.08	0.58
32:B6:42:TRP:HA	32:B6:42:TRP:HE3	1.69	0.58
33:B7:7:PRO:CB	35:BA:1309:G:H4'	2.34	0.58
34:B8:56:GLU:HA	34:B8:59:LYS:NZ	2.18	0.58
35:BA:1049:C:H2'	35:BA:1050:A:C8	2.38	0.58
35:BA:1469:A:O2'	35:BA:1470:G:H5'	2.03	0.58
35:BA:15:G:H2'	35:BA:16:G:C8	2.34	0.58
35:BA:1710:C:H2'	35:BA:1711:C:C6	2.39	0.58
35:BA:1789:A:H2'	35:BA:1790:C:H6	1.67	0.58
35:BA:1997:G:O2'	35:BA:1998:G:H5'	2.03	0.58
35:BA:1998:G:H2'	35:BA:1999:C:H6	1.69	0.58
35:BA:2081:C:O2'	35:BA:2082:A:H5'	2.04	0.58
35:BA:2538:C:C2'	35:BA:2539:C:H5'	2.32	0.58
35:BA:272:G:H1'	35:BA:272(B):G:O4'	2.03	0.58
35:BA:493:G:H2'	35:BA:494:G:H5''	1.85	0.58
35:BA:668:G:C2	35:BA:670:A:C6	2.92	0.58
35:BA:76:C:H2'	35:BA:77:C:H6	1.68	0.58
35:BA:782:A:C2	38:BD:226:MET:HE2	2.39	0.58
36:BB:83:G:O2'	36:BB:84:C:H5'	2.04	0.58
38:BD:172:TYR:HD1	38:BD:186:HIS:CA	2.16	0.58
38:BD:35:LYS:HD3	38:BD:63:ARG:CB	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:1:MET:N	39:BE:84:PHE:HB2	2.18	0.58
41:BG:115:ARG:HH22	41:BG:136:ARG:CD	2.15	0.58
41:BG:138:GLN:CD	41:BG:153:ARG:H	2.07	0.58
45:BO:62:VAL:HG12	45:BO:63:VAL:N	2.18	0.58
50:BT:110:ILE:HA	50:BT:113:LYS:HD2	1.85	0.58
50:BT:13:ARG:HH12	50:BT:15:VAL:HG13	1.64	0.58
1:CA:1294:G:H2'	1:CA:1295:G:H8	1.69	0.58
1:CA:1341:U:O2'	1:CA:1342:C:H5'	2.03	0.58
1:CA:1365:G:O2'	1:CA:1366:C:H5'	2.03	0.58
1:CA:304:U:H2'	1:CA:305:G:C8	2.39	0.58
1:CA:781:A:H2'	1:CA:782:A:H5'	1.84	0.58
2:CB:67:THR:CG2	2:CB:155:LEU:HG	2.33	0.58
2:CB:163:PHE:HA	2:CB:185:ILE:O	2.03	0.58
2:CB:50:GLU:OE1	2:CB:200:ILE:HB	2.03	0.58
3:CC:91:LEU:O	3:CC:94:LEU:HG	2.03	0.58
4:CD:64:LEU:HB2	4:CD:198:VAL:HG11	1.86	0.58
6:CF:37:VAL:HG13	6:CF:65:VAL:CG1	2.33	0.58
8:CH:29:SER:O	8:CH:32:LYS:HB2	2.03	0.58
12:CL:5:PRO:HG2	12:CL:10:LEU:HD21	1.86	0.58
12:CL:119:LYS:O	12:CL:120:TYR:HB2	2.03	0.58
12:CL:119:LYS:C	12:CL:120:TYR:HD1	2.05	0.58
16:CP:68:ASP:O	16:CP:71:ARG:HB3	2.02	0.58
18:CR:22:VAL:HA	18:CR:25:THR:OG1	2.03	0.58
21:CU:2:GLY:N	21:CU:5:ASP:HB2	2.18	0.58
23:CW:17:C:H4'	23:CW:62:C:C5'	2.33	0.58
27:D1:73:LEU:HD23	27:D1:90:ILE:HG22	1.86	0.58
33:D7:31:LEU:CD2	33:D7:42:LEU:HB3	2.32	0.58
35:DA:1049:C:H2'	35:DA:1050:A:C8	2.38	0.58
35:DA:1108:U:C2'	35:DA:1109:C:H5'	2.33	0.58
35:DA:1255:U:H5''	35:DA:1256:G:H5''	1.84	0.58
35:DA:2571:C:H5'	35:DA:2572:A:H5'	1.86	0.58
35:DA:2707:G:H2'	35:DA:2708:G:H8	1.68	0.58
35:DA:401:A:H2'	35:DA:402:A:C8	2.38	0.58
35:DA:662:G:H2'	35:DA:663:G:H8	1.68	0.58
35:DA:861:A:H62	35:DA:916:G:H21	1.49	0.58
35:DA:926:A:H8	35:DA:926:A:H5'	1.67	0.58
36:DB:78:A:C2	36:DB:100:A:C4	2.91	0.58
38:DD:134:ARG:O	38:DD:136:ILE:N	2.36	0.58
38:DD:172:TYR:HD1	38:DD:186:HIS:CA	2.16	0.58
39:DE:34:VAL:CG2	39:DE:48:GLN:HE21	2.16	0.58
39:DE:6:GLY:HA2	39:DE:51:PHE:HE2	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:125:PHE:HE2	41:DG:173:LEU:HD12	1.69	0.58
41:DG:38:VAL:HG13	41:DG:91:ARG:HD3	1.86	0.58
42:DH:85:LYS:HE3	42:DH:144:VAL:HB	1.83	0.58
43:DI:88:ILE:CG2	43:DI:89:TYR:H	2.16	0.58
44:DN:23:LEU:HB3	44:DN:60:ILE:HG21	1.84	0.58
45:DO:13:ASN:HD21	45:DO:96:THR:H	1.50	0.58
45:DO:14:THR:HG22	45:DO:52:VAL:HG21	1.85	0.58
45:DO:31:LYS:HB3	45:DO:32:TYR:CD1	2.37	0.58
40:DF:34:TRP:CB	46:DP:11:GLY:HA3	2.33	0.58
47:DQ:9:TYR:O	47:DQ:10:ARG:CG	2.44	0.58
50:DT:13:ARG:HH12	50:DT:15:VAL:HG13	1.68	0.58
51:DU:3:ARG:HG3	51:DU:3:ARG:O	2.02	0.58
52:DV:34:GLU:HG2	52:DV:35:LEU:N	2.18	0.58
55:DY:10:GLY:HA2	55:DY:27:VAL:CG1	2.24	0.58
55:DY:28:LYS:HE3	55:DY:28:LYS:O	2.03	0.58
56:DZ:22:GLY:O	56:DZ:41:LEU:HD21	2.03	0.58
56:DZ:48:PHE:CA	56:DZ:51:ALA:HB3	2.33	0.58
1:AA:1522:U:C2'	1:AA:1523:G:H5'	2.34	0.58
1:AA:506:G:H2'	1:AA:507:C:C6	2.39	0.58
1:AA:661:G:H2'	1:AA:662:G:H8	1.67	0.58
1:AA:735:C:H2'	1:AA:736:C:C6	2.37	0.58
4:AD:64:LEU:HB2	4:AD:198:VAL:HG11	1.85	0.58
6:AF:45:LEU:O	6:AF:46:ARG:HD2	2.03	0.58
7:AG:146:GLU:O	7:AG:149:ARG:HB2	2.03	0.58
5:AE:78:HIS:HD2	8:AH:104:ARG:NE	2.01	0.58
11:AK:69:ALA:O	11:AK:73:MET:N	2.29	0.58
14:AN:40:CYS:SG	14:AN:41:ARG:N	2.77	0.58
20:AT:50:GLU:HG3	20:AT:51:GLU:H	1.69	0.58
23:AW:29:C:H2'	23:AW:30:G:H8	1.68	0.58
25:AY:34:ASN:CG	25:AY:35:PRO:HD2	2.24	0.58
28:B2:53:LEU:CA	28:B2:56:GLN:HG2	2.30	0.58
32:B6:51:GLU:CG	32:B6:52:VAL:H	2.00	0.58
35:BA:1034:G:H22	35:BA:1122:G:H1'	1.66	0.58
35:BA:1707:G:H2'	35:BA:1708:C:C6	2.38	0.58
35:BA:214:G:O2'	35:BA:215:G:O4'	2.20	0.58
35:BA:2863:C:C2'	35:BA:2864:G:H5''	2.34	0.58
35:BA:611:C:O2'	35:BA:612:C:H5'	2.03	0.58
35:BA:820:A:H2'	35:BA:821:A:O4'	2.03	0.58
35:BA:841:A:H2'	35:BA:842:G:C8	2.39	0.58
37:BC:76:ALA:HB3	37:BC:94:VAL:HG11	1.85	0.58
38:BD:69:ARG:HD3	38:BD:105:ILE:HD12	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:34:VAL:CG2	39:BE:48:GLN:HE21	2.15	0.58
40:BF:110:LEU:HD21	40:BF:181:LEU:HD23	1.85	0.58
43:BI:88:ILE:HG23	43:BI:89:TYR:H	1.68	0.58
44:BN:57:ALA:HB1	44:BN:60:ILE:HD11	1.84	0.58
46:BP:62:LEU:HD22	46:BP:62:LEU:O	2.03	0.58
46:BP:91:PHE:N	46:BP:91:PHE:CD1	2.71	0.58
46:BP:95:VAL:O	46:BP:125:VAL:HG23	2.03	0.58
35:BA:2876:G:H4'	50:BT:2:ASN:O	2.03	0.58
51:BU:49:HIS:HA	51:BU:52:ARG:HB2	1.86	0.58
54:BX:65:ARG:CA	54:BX:65:ARG:NE	2.66	0.58
54:BX:59:VAL:HG23	54:BX:74:PRO:CD	2.33	0.58
55:BY:68:HIS:ND1	55:BY:69:ALA:N	2.51	0.58
55:BY:95:LYS:HG2	55:BY:100:ALA:CA	2.18	0.58
56:BZ:116:VAL:HG12	56:BZ:117:LEU:N	2.15	0.58
56:BZ:97:GLU:HA	56:BZ:126:VAL:O	2.03	0.58
1:CA:1515:C:O2'	1:CA:1516:G:H5'	2.03	0.58
1:CA:178:C:H2'	1:CA:179:A:H8	1.68	0.58
2:CB:167:PRO:HD2	2:CB:188:ALA:HB2	1.85	0.58
2:CB:178:ARG:HH22	2:CB:196:LEU:HA	1.68	0.58
2:CB:185:ILE:HG22	2:CB:199:TYR:HD1	1.68	0.58
2:CB:71:VAL:CG2	2:CB:93:VAL:HG23	2.33	0.58
6:CF:69:GLU:HG2	6:CF:70:ASP:H	1.68	0.58
12:CL:32:PHE:HB3	12:CL:84:LEU:HD21	1.86	0.58
12:CL:86:ARG:CG	12:CL:87:GLY:H	2.16	0.58
15:CO:50:HIS:O	15:CO:53:HIS:HB3	2.02	0.58
33:D7:8:ASN:C	33:D7:8:ASN:HD22	2.06	0.58
35:DA:1034:G:H22	35:DA:1122:G:H1'	1.66	0.58
35:DA:1198:U:H2'	35:DA:1199:U:C6	2.38	0.58
35:DA:1224:C:O3'	52:DV:88:ARG:HB3	2.03	0.58
35:DA:1324:G:H3'	35:DA:1325:G:C5'	2.33	0.58
35:DA:1503:U:H2'	35:DA:1504:C:C5	2.37	0.58
35:DA:195:A:H5''	35:DA:196:A:OP2	2.03	0.58
35:DA:2078:C:O2'	35:DA:2079:U:H5'	2.04	0.58
35:DA:2600:A:HO2'	35:DA:2601:C:H5'	1.67	0.58
35:DA:2840:C:H2'	35:DA:2841:C:H6	1.69	0.58
35:DA:493:G:H3'	35:DA:494:G:H5''	1.85	0.58
35:DA:580:C:H2'	35:DA:581:C:C6	2.38	0.58
35:DA:680:G:H2'	35:DA:681:G:H8	1.68	0.58
38:DD:35:LYS:HE2	38:DD:104:TYR:CB	2.33	0.58
39:DE:24:THR:OG1	39:DE:188:VAL:HG11	2.04	0.58
44:DN:110:GLY:HA2	44:DN:114:ARG:NH2	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:55:VAL:HG12	44:DN:126:PRO:HA	1.84	0.58
55:DY:37:VAL:O	55:DY:38:ILE:HB	2.03	0.58
55:DY:42:VAL:HB	55:DY:65:ALA:HB3	1.83	0.58
1:AA:320:C:O2'	1:AA:1435:G:H1'	2.03	0.58
1:AA:726:C:H2'	1:AA:727:G:C8	2.39	0.58
1:AA:769:G:H2'	1:AA:770:C:H6	1.67	0.58
1:AA:779:C:O2'	1:AA:780:A:H5'	2.02	0.58
1:AA:794:A:H2'	1:AA:795:C:C6	2.38	0.58
2:AB:73:THR:HG22	2:AB:93:VAL:O	2.03	0.58
4:AD:11:LEU:O	4:AD:13:ARG:N	2.36	0.58
6:AF:82:ARG:HB2	6:AF:85:VAL:CG2	2.34	0.58
9:AI:26:VAL:HG22	9:AI:61:ALA:HB3	1.86	0.58
10:AJ:50:ILE:N	10:AJ:50:ILE:HD13	2.01	0.58
19:AS:12:ASP:O	19:AS:16:LEU:HD13	2.03	0.58
20:AT:17:ARG:HA	20:AT:20:LEU:HD12	1.84	0.58
20:AT:73:HIS:HB3	20:AT:74:LYS:CD	2.31	0.58
23:AW:23:G:H2'	23:AW:24:C:H5''	1.86	0.58
25:AY:13:HIS:O	25:AY:132:ILE:HD13	2.03	0.58
25:AY:58:VAL:HG22	25:AY:68:VAL:HG22	1.84	0.58
27:B1:87:PRO:HB2	27:B1:91:LYS:HZ1	1.68	0.58
34:B8:49:VAL:CG1	34:B8:53:PRO:HD3	2.33	0.58
35:BA:1259:G:O2'	35:BA:1260:G:H5'	2.03	0.58
35:BA:127:A:H5''	35:BA:128:C:O4'	2.04	0.58
35:BA:1447:G:H2'	35:BA:1448:G:H8	1.68	0.58
35:BA:1813:G:H1'	38:BD:50:THR:OG1	2.04	0.58
35:BA:1854:A:H3'	35:BA:1855:G:H8	1.68	0.58
35:BA:2033:A:H4'	35:BA:2034:U:OP1	2.03	0.58
35:BA:2555:U:H2'	35:BA:2556:C:H5'	1.84	0.58
35:BA:271(Q):G:H2'	35:BA:271(R):G:H8	1.68	0.58
35:BA:2808:U:O2'	35:BA:2809:A:H5'	2.03	0.58
35:BA:2870:C:C2'	35:BA:2871:C:H5'	2.34	0.58
39:BE:197:ILE:HD11	39:BE:199:ARG:HH22	1.67	0.58
39:BE:51:PHE:H	39:BE:74:PRO:CB	2.16	0.58
39:BE:59:VAL:HG21	39:BE:63:LEU:HA	1.85	0.58
40:BF:20:LEU:HB3	40:BF:23:ASP:OD2	2.04	0.58
41:BG:91:ARG:HG2	41:BG:92:VAL:N	2.18	0.58
42:BH:109:PHE:CE1	42:BH:152:ARG:NE	2.71	0.58
46:BP:90:ARG:CD	46:BP:91:PHE:HD1	2.17	0.58
48:BR:36:THR:HB	48:BR:40:LYS:HD2	1.86	0.58
49:BS:85:VAL:CG2	49:BS:106:ARG:HB2	2.34	0.58
50:BT:74:ARG:HG2	50:BT:74:ARG:HH11	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:BX:35:THR:HB	54:BX:75:ASP:OD2	2.03	0.58
1:CA:1097:C:O2'	1:CA:1098:C:H5'	2.02	0.58
1:CA:1292:U:O2'	1:CA:1293:G:H5'	2.02	0.58
1:CA:779:C:O2'	1:CA:780:A:H5'	2.02	0.58
2:CB:102:LEU:CD1	2:CB:102:LEU:N	2.65	0.58
2:CB:19:HIS:CD2	2:CB:20:GLU:HG2	2.38	0.58
4:CD:127:THR:HG23	4:CD:147:ALA:HB3	1.86	0.58
6:CF:75:LEU:HD22	6:CF:79:LEU:HD11	1.85	0.58
10:CJ:32:ALA:N	10:CJ:78:ASN:ND2	2.51	0.58
14:CN:51:GLY:C	14:CN:53:LEU:H	2.07	0.58
20:CT:42:GLN:O	20:CT:43:LEU:HD23	2.03	0.58
25:CY:180:GLU:O	25:CY:183:ILE:HG13	2.03	0.58
27:D1:39:LYS:HE3	35:DA:201:C:P	2.44	0.58
35:DA:1771:C:H1'	35:DA:1786:A:C8	2.38	0.58
35:DA:1792:G:O2'	35:DA:1793:C:H5'	2.03	0.58
35:DA:2334:G:H5'	49:DS:13:ARG:CG	2.29	0.58
35:DA:2825:C:H2'	35:DA:2826:A:O4'	2.03	0.58
35:DA:845:G:O2'	35:DA:846:C:H5	1.86	0.58
39:DE:81:ILE:O	39:DE:81:ILE:HG22	2.04	0.58
43:DI:66:GLU:O	43:DI:70:GLU:HG2	2.04	0.58
44:DN:129:PRO:O	44:DN:130:HIS:HB3	2.03	0.58
44:DN:26:LEU:HD21	44:DN:99:LEU:HD11	1.84	0.58
35:DA:2563:U:O2'	45:DO:28:SER:HB3	2.02	0.58
35:DA:1190:G:C5'	46:DP:35:HIS:HA	2.34	0.58
51:DU:40:PHE:CD2	51:DU:40:PHE:N	2.69	0.58
54:DX:49:VAL:CG1	54:DX:50:LYS:H	2.13	0.58
55:DY:28:LYS:CE	55:DY:30:VAL:HA	2.32	0.58
1:AA:1248:A:H2'	1:AA:1249:C:H5'	1.84	0.58
1:AA:708:C:O2'	1:AA:709:G:H5'	2.02	0.58
1:AA:802:A:H2'	1:AA:803:G:C5'	2.34	0.58
1:AA:877:C:H5''	8:AH:88:LYS:HD2	1.85	0.58
2:AB:101:MET:O	2:AB:108:ILE:HG21	2.03	0.58
2:AB:71:VAL:O	2:AB:164:VAL:HG22	2.04	0.58
4:AD:133:VAL:HG12	4:AD:135:LEU:H	1.69	0.58
5:AE:6:PHE:HB3	5:AE:35:GLY:O	2.03	0.58
5:AE:78:HIS:HB2	5:AE:79:GLU:OE1	2.03	0.58
8:AH:103:VAL:HG12	8:AH:108:GLY:HA3	1.83	0.58
18:AR:74:ARG:HH11	18:AR:74:ARG:HG3	1.69	0.58
20:AT:16:HIS:O	20:AT:19:SER:HB2	2.04	0.58
25:AY:171:LYS:HA	25:AY:174:GLN:NE2	2.18	0.58
34:B8:60:LEU:HA	34:B8:63:PRO:CG	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1006:C:H2'	35:BA:1007:C:C6	2.38	0.58
35:BA:1150:C:C2'	35:BA:1151:G:H5'	2.34	0.58
35:BA:1677:A:H2'	35:BA:1678:G:H8	1.67	0.58
35:BA:2098:U:H2'	35:BA:2099:U:C6	2.39	0.58
35:BA:221:A:H4'	35:BA:222:A:O5'	2.04	0.58
35:BA:2314:C:H2'	35:BA:2315:G:H8	1.69	0.58
35:BA:2368:C:H2'	35:BA:2369:A:H8	1.67	0.58
35:BA:256:A:H2'	35:BA:257:A:C8	2.39	0.58
35:BA:2821:A:H3'	35:BA:2821:A:OP2	2.03	0.58
35:BA:470:A:OP1	40:BF:59:TYR:HE2	1.87	0.58
40:BF:31:HIS:O	40:BF:32:LEU:C	2.42	0.58
45:BO:104:ARG:HH12	50:BT:35:LYS:HD3	1.68	0.58
47:BQ:35:VAL:CG1	47:BQ:130:LYS:HB3	2.34	0.58
47:BQ:43:THR:OG1	47:BQ:45:GLN:HB2	2.03	0.58
47:BQ:81:VAL:HG23	47:BQ:82:ARG:HH11	1.69	0.58
48:BR:12:ARG:O	48:BR:13:HIS:HB3	2.04	0.58
35:BA:1275:A:C5	48:BR:16:HIS:ND1	2.72	0.58
49:BS:23:ARG:O	49:BS:86:ALA:HB3	2.04	0.58
53:BW:110:LYS:HG3	53:BW:111:HIS:ND1	2.19	0.58
54:BX:88:LYS:C	54:BX:90:GLU:H	2.07	0.58
56:BZ:116:VAL:HG12	56:BZ:117:LEU:HD23	1.86	0.58
1:CA:936:C:O2'	1:CA:937:A:H5'	2.02	0.58
3:CC:107:GLN:O	3:CC:108:ASN:HB2	2.01	0.58
6:CF:43:LEU:H	6:CF:43:LEU:CD1	2.17	0.58
7:CG:100:ALA:C	7:CG:104:LEU:HD23	2.24	0.58
8:CH:11:THR:HA	8:CH:14:ARG:CZ	2.33	0.58
10:CJ:6:ILE:HD12	10:CJ:6:ILE:O	2.03	0.58
13:CM:23:TYR:OH	13:CM:71:ARG:HD3	2.04	0.58
14:CN:23:ARG:HD3	14:CN:29:ARG:O	2.04	0.58
16:CP:73:LEU:O	16:CP:77:ALA:HB2	2.03	0.58
17:CQ:86:GLU:C	17:CQ:88:TYR:N	2.57	0.58
20:CT:76:ALA:O	20:CT:80:ARG:HG2	2.04	0.58
23:CW:36:A:O2'	23:CW:37:U:H5'	2.03	0.58
25:CY:160:GLU:O	25:CY:161:ILE:C	2.41	0.58
25:CY:32:ARG:NE	25:CY:32:ARG:CA	2.66	0.58
26:D0:1:MET:O	26:D0:2:ALA:HB3	2.04	0.58
35:DA:1596:A:H5'	35:DA:1597:A:OP2	2.04	0.58
35:DA:1615:C:H5	35:DA:1617:C:C4	2.21	0.58
35:DA:2012:G:OP1	53:DW:98:LYS:HA	2.03	0.58
35:DA:2052:G:C2	39:DE:149:ARG:HA	2.38	0.58
35:DA:2264:C:H2'	35:DA:2265:U:C6	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:7:G:H4'	44:DN:13:TRP:HH2	1.68	0.58
35:DA:92:A:H2'	35:DA:93:G:C8	2.39	0.58
35:DA:972:G:H2'	35:DA:973:A:C8	2.39	0.58
40:DF:67:GLN:O	40:DF:67:GLN:CG	2.35	0.58
44:DN:137:LYS:HG2	44:DN:138:LEU:N	2.17	0.58
46:DP:107:LYS:C	46:DP:109:GLY:H	2.07	0.58
48:DR:37:THR:HA	48:DR:111:LEU:HA	1.85	0.58
49:DS:84:GLN:HA	49:DS:105:ALA:O	2.03	0.58
49:DS:23:ARG:O	49:DS:86:ALA:HB3	2.03	0.58
52:DV:72:VAL:HG12	52:DV:73:SER:N	2.11	0.58
54:DX:60:ARG:HG3	54:DX:71:GLY:HA3	1.86	0.58
1:AA:112:G:O2'	1:AA:113:G:H5'	2.04	0.58
1:AA:361:G:O2'	1:AA:362:G:H5'	2.03	0.58
1:AA:788:U:C4	1:AA:789:U:C5	2.92	0.58
1:AA:936:C:O2'	1:AA:937:A:H5'	2.03	0.58
3:AC:88:ARG:HG2	3:AC:101:LEU:HB2	1.86	0.58
4:AD:100:ARG:CZ	4:AD:137:SER:HA	2.33	0.58
4:AD:18:LYS:CE	4:AD:31:CYS:HB3	2.33	0.58
10:AJ:23:ILE:HG22	10:AJ:23:ILE:O	2.04	0.58
10:AJ:75:ILE:HG13	10:AJ:76:ASN:N	2.18	0.58
12:AL:70:ILE:HG13	12:AL:100:ILE:HD12	1.85	0.58
1:AA:127:G:HO2'	17:AQ:2:PRO:N	2.01	0.58
25:AY:164:ILE:HG22	25:AY:165:THR:N	2.19	0.58
26:B0:38:VAL:HG23	26:B0:59:LEU:HB2	1.85	0.58
28:B2:31:GLU:O	28:B2:37:PHE:HB2	2.03	0.58
33:B7:19:ARG:HD3	35:BA:125:G:H5''	1.84	0.58
35:BA:2707:G:H2'	35:BA:2708:G:H8	1.68	0.58
35:BA:268:C:H42	35:BA:424:G:H1	1.52	0.58
35:BA:42:G:H2'	35:BA:43:A:C8	2.39	0.58
35:BA:66:C:H2'	35:BA:67:U:H5'	1.86	0.58
38:BD:48:ARG:NH1	38:BD:48:ARG:HG3	2.18	0.58
40:BF:3:GLU:HA	40:BF:24:LEU:HB3	1.85	0.58
41:BG:142:PRO:HG2	41:BG:143:GLU:H	1.69	0.58
44:BN:78:TYR:N	44:BN:79:PRO:HD2	2.18	0.58
46:BP:144:GLU:N	46:BP:145:PRO:HD3	2.18	0.58
50:BT:101:PHE:CD2	50:BT:102:ILE:N	2.70	0.58
50:BT:35:LYS:HZ1	50:BT:41:ARG:HE	1.52	0.58
52:BV:18:LEU:CD2	52:BV:19:LYS:H	2.08	0.58
52:BV:34:GLU:CB	52:BV:62:LEU:HD12	2.33	0.58
55:BY:98:VAL:O	55:BY:99:CYS:SG	2.62	0.58
1:CA:10:A:H2'	1:CA:11:G:H8	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:192:U:H2'	1:CA:193:C:H6	1.68	0.58
1:CA:337:C:H2'	1:CA:338:A:C8	2.35	0.58
1:CA:569:C:H42	1:CA:881:G:H1	1.51	0.58
1:CA:680:C:O2'	1:CA:681:C:H5'	2.04	0.58
1:CA:758:G:H8	1:CA:758:G:O5'	1.86	0.58
2:CB:224:GLN:O	2:CB:224:GLN:HG2	2.02	0.58
2:CB:79:ASP:O	2:CB:82:ARG:N	2.36	0.58
3:CC:130:VAL:HA	3:CC:133:ALA:HB3	1.86	0.58
3:CC:130:VAL:O	3:CC:134:ILE:HG13	2.04	0.58
6:CF:10:LEU:HD13	6:CF:61:LEU:HD11	1.86	0.58
7:CG:146:GLU:O	7:CG:149:ARG:HB2	2.03	0.58
15:CO:70:LEU:HD23	15:CO:78:TYR:HA	1.85	0.58
16:CP:53:VAL:HG23	16:CP:54:GLU:H	1.68	0.58
16:CP:71:ARG:HA	16:CP:74:LEU:HD12	1.86	0.58
27:D1:15:ALA:O	27:D1:46:LEU:HD23	2.02	0.58
32:D6:10:LEU:HD22	32:D6:10:LEU:N	2.18	0.58
35:DA:1448:G:H2'	35:DA:1449:A:C8	2.38	0.58
35:DA:2114:A:C2'	35:DA:2115:G:H5'	2.34	0.58
35:DA:2241:A:H2'	35:DA:2242:G:H8	1.69	0.58
35:DA:2863:C:C2'	35:DA:2864:G:H5''	2.34	0.58
35:DA:470:A:C2	35:DA:471:A:C4	2.91	0.58
37:DC:196:LEU:C	37:DC:198:ALA:H	2.07	0.58
38:DD:211:ARG:CG	38:DD:211:ARG:HH11	2.17	0.58
38:DD:83:GLU:HB2	38:DD:92:ILE:HD11	1.86	0.58
35:DA:2619:C:OP1	39:DE:152:LYS:HD3	2.04	0.58
39:DE:49:LEU:O	39:DE:78:LEU:HA	2.03	0.58
41:DG:38:VAL:CA	41:DG:158:ALA:HB3	2.34	0.58
42:DH:136:ILE:HD12	42:DH:136:ILE:N	2.17	0.58
42:DH:89:ILE:CD1	42:DH:129:THR:HB	2.33	0.58
43:DI:8:PRO:HD3	43:DI:15:VAL:CG1	2.34	0.58
44:DN:24:GLY:HA2	44:DN:27:ALA:HB3	1.84	0.58
46:DP:62:LEU:O	46:DP:62:LEU:HD22	2.03	0.58
46:DP:85:LEU:HA	46:DP:88:LEU:CB	2.30	0.58
49:DS:87:PHE:O	49:DS:88:ASP:HB2	2.02	0.58
50:DT:129:ARG:CZ	50:DT:131:ALA:HB3	2.33	0.58
47:DQ:141:GLN:C	56:DZ:53:ILE:HB	2.24	0.58
1:AA:1223:C:P	1:AA:1224:G:H2'	2.43	0.58
1:AA:1435:G:H2'	1:AA:1436:U:C6	2.39	0.58
1:AA:1479:C:H2'	1:AA:1480:G:C8	2.39	0.58
1:AA:441:A:H3'	1:AA:442:C:C6	2.38	0.58
1:AA:460:G:O6	1:AA:470:C:H5''	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:881:G:P	12:AL:12:ARG:HH22	2.27	0.58
4:AD:14:ARG:O	4:AD:16:GLY:N	2.36	0.58
7:AG:29:LYS:HB2	7:AG:105:VAL:HG21	1.84	0.58
7:AG:121:ALA:CA	7:AG:124:LEU:HD12	2.34	0.58
7:AG:143:ARG:O	7:AG:147:ALA:HB2	2.03	0.58
8:AH:11:THR:HA	8:AH:14:ARG:HH12	1.68	0.58
9:AI:79:LEU:CD2	9:AI:102:LEU:HA	2.32	0.58
9:AI:4:TYR:HB2	9:AI:19:LEU:HD12	1.86	0.58
1:AA:255:G:O3'	17:AQ:17:LYS:HD2	2.02	0.58
19:AS:14:HIS:CD2	19:AS:15:LEU:HD22	2.39	0.58
25:AY:170:ALA:C	25:AY:172:ALA:N	2.58	0.58
27:B1:87:PRO:O	27:B1:89:GLU:N	2.37	0.58
35:BA:1337:G:H2'	35:BA:1338:G:O4'	2.04	0.58
35:BA:1357:U:O2'	35:BA:1358:G:H5'	2.03	0.58
35:BA:2302:G:H1'	41:BG:128:ARG:HG3	1.86	0.58
35:BA:332:A:H4'	35:BA:333:G:OP1	2.03	0.58
35:BA:45:C:H2'	35:BA:47:C:H6	1.68	0.58
35:BA:61:G:H1	35:BA:94:C:H42	1.52	0.58
35:BA:94(A):G:C3'	35:BA:95:G:H5''	2.33	0.58
38:BD:10:THR:O	38:BD:11:PRO:C	2.38	0.58
38:BD:134:ARG:HB2	38:BD:135:PHE:HD1	1.69	0.58
39:BE:49:LEU:O	39:BE:78:LEU:HA	2.03	0.58
42:BH:136:ILE:HD12	42:BH:136:ILE:N	2.15	0.58
42:BH:20:ALA:HB1	42:BH:21:PRO:CD	2.34	0.58
42:BH:40:GLU:HB2	42:BH:41:MET:SD	2.43	0.58
43:BI:66:GLU:O	43:BI:70:GLU:HG2	2.04	0.58
43:BI:94:ALA:C	43:BI:96:ASP:N	2.57	0.58
44:BN:107:LEU:HD12	44:BN:108:PRO:O	2.03	0.58
48:BR:13:HIS:CE1	48:BR:16:HIS:HB2	2.39	0.58
48:BR:6:SER:HA	48:BR:8:ARG:NH2	2.19	0.58
52:BV:22:VAL:HG21	52:BV:96:ILE:HD12	1.86	0.58
52:BV:38:LEU:CD2	52:BV:39:LEU:N	2.65	0.58
52:BV:75:PHE:HD1	52:BV:87:HIS:O	1.87	0.58
53:BW:1:MET:HG3	53:BW:2:GLU:N	2.19	0.58
28:B2:29:LYS:HZ1	54:BX:9:LEU:HA	1.67	0.58
47:BQ:134:ARG:HH12	56:BZ:119:GLU:CD	2.06	0.58
1:CA:129(A):G:N2	1:CA:189(E):U:H1'	2.19	0.58
1:CA:453:A:C5	1:CA:454:C:C4	2.92	0.58
1:CA:826:C:H2'	1:CA:827:U:C6	2.39	0.58
1:CA:995:C:O2'	1:CA:996:A:H5'	2.04	0.58
2:CB:222:ILE:HG23	2:CB:223:ILE:N	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:69:LEU:CB	2:CB:162:ILE:HG22	2.33	0.58
4:CD:14:ARG:O	4:CD:16:GLY:N	2.36	0.58
4:CD:176:LEU:CD1	4:CD:177:ASP:H	2.16	0.58
7:CG:29:LYS:HB2	7:CG:105:VAL:HG21	1.86	0.58
11:CK:62:GLN:C	11:CK:64:ALA:H	2.07	0.58
12:CL:9:GLN:O	12:CL:11:VAL:N	2.36	0.58
13:CM:78:ILE:HA	13:CM:81:LEU:CD1	2.33	0.58
1:CA:127:G:HO2'	17:CQ:2:PRO:N	2.02	0.58
25:CY:162:GLN:HG3	25:CY:166:ASP:OD2	2.04	0.58
27:D1:13:ILE:CD1	27:D1:14:VAL:HG12	2.34	0.58
31:D5:2:ALA:N	35:DA:747:U:N3	2.51	0.58
31:D5:8:LYS:O	31:D5:9:LYS:HD2	2.03	0.58
35:DA:1822:G:H2'	35:DA:1823:G:H8	1.68	0.58
35:DA:2410:G:N2	35:DA:2411:A:H1'	2.19	0.58
35:DA:2468:G:HO2'	35:DA:2476:A:H8	1.52	0.58
35:DA:8:A:H2'	35:DA:9:U:C6	2.37	0.58
38:DD:181:GLU:HA	38:DD:272:ALA:O	2.03	0.58
40:DF:3:GLU:CB	40:DF:24:LEU:HG	2.33	0.58
41:DG:45:GLU:O	41:DG:47:LYS:N	2.36	0.58
45:DO:11:ALA:HB1	45:DO:99:PHE:O	2.04	0.58
46:DP:21:ARG:HH11	46:DP:21:ARG:HG3	1.68	0.58
47:DQ:114:ALA:C	47:DQ:116:GLU:H	2.05	0.58
50:DT:50:ILE:N	50:DT:50:ILE:HD12	2.19	0.58
54:DX:59:VAL:HG23	54:DX:74:PRO:CD	2.32	0.58
55:DY:76:CYS:SG	55:DY:77:PRO:CD	2.92	0.58
1:AA:9:G:H2'	1:AA:10:A:C8	2.39	0.58
1:AA:1202:G:C2'	1:AA:1203:C:H5'	2.34	0.58
1:AA:680:C:O2'	1:AA:681:C:H5'	2.04	0.58
2:AB:158:LEU:H	2:AB:158:LEU:HD12	1.68	0.58
2:AB:55:PHE:CD1	2:AB:58:ILE:HD12	2.39	0.58
2:AB:80:ILE:HG13	2:AB:81:VAL:N	2.18	0.58
7:AG:152:ALA:O	7:AG:154:TYR:N	2.36	0.58
7:AG:75:VAL:HA	7:AG:88:PRO:HA	1.86	0.58
8:AH:27:PRO:HA	8:AH:58:TYR:HA	1.86	0.58
9:AI:5:TYR:HA	9:AI:17:VAL:O	2.03	0.58
9:AI:17:VAL:CG2	9:AI:81:ILE:HD13	2.34	0.58
11:AK:65:ALA:O	11:AK:68:ALA:HB3	2.04	0.58
1:AA:528:C:N4	12:AL:49:ASN:HD22	2.01	0.58
13:AM:9:ILE:CG2	13:AM:11:ARG:HG3	2.33	0.58
1:AA:656:C:H4'	15:AO:62:GLN:HE22	1.68	0.58
15:AO:82:ILE:O	15:AO:82:ILE:HD13	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:65:ILE:N	17:AQ:65:ILE:HD12	2.19	0.58
19:AS:49:ILE:O	19:AS:60:VAL:HG12	2.04	0.58
34:B8:60:LEU:CD2	34:B8:60:LEU:N	2.67	0.58
35:BA:1024:G:C8	35:BA:1025:G:H2'	2.39	0.58
35:BA:1427:A:H4'	35:BA:1428:C:O5'	2.04	0.58
35:BA:1744:C:O2'	35:BA:1745:C:H5'	2.03	0.58
32:B6:27:LYS:HD2	35:BA:2285:C:OP2	2.04	0.58
35:BA:360:G:H2'	35:BA:361:G:H8	1.69	0.58
35:BA:40:C:H2'	35:BA:41:C:H6	1.68	0.58
35:BA:635:C:O2'	35:BA:636:G:H5'	2.04	0.58
35:BA:977:G:O2'	35:BA:978:G:H5'	2.04	0.58
38:BD:35:LYS:HE2	38:BD:104:TYR:CG	2.39	0.58
35:BA:1824:G:OP1	38:BD:52:ARG:HD3	2.04	0.58
39:BE:78:LEU:CD2	39:BE:78:LEU:N	2.67	0.58
45:BO:63:VAL:HG11	45:BO:85:VAL:HG23	1.86	0.58
46:BP:106:LEU:HD11	46:BP:112:LEU:HB2	1.86	0.58
51:BU:110:VAL:O	51:BU:114:LYS:N	2.31	0.58
51:BU:47:TYR:HA	51:BU:50:ARG:NH1	2.18	0.58
56:BZ:10:ARG:HB3	56:BZ:36:LYS:CB	2.29	0.58
1:CA:1396:A:O4'	1:CA:1398:A:H1'	2.03	0.58
1:CA:524:G:H2'	1:CA:525:C:C6	2.38	0.58
1:CA:539:A:OP1	12:CL:114:LYS:HE2	2.04	0.58
1:CA:601:C:O2'	1:CA:602:A:H5'	2.04	0.58
7:CG:121:ALA:HA	7:CG:124:LEU:HD12	1.84	0.58
1:CA:823:G:H21	8:CH:1:MET:HE1	1.69	0.58
1:CA:796:C:P	11:CK:123:LYS:HZ2	2.27	0.58
12:CL:83:VAL:CG2	12:CL:84:LEU:H	2.08	0.58
17:CQ:40:LYS:HG2	17:CQ:41:LYS:N	2.19	0.58
18:CR:56:THR:HG21	18:CR:63:GLN:HE22	1.69	0.58
19:CS:41:VAL:CB	19:CS:44:MET:HB2	2.34	0.58
22:CV:27:G:O2'	22:CV:28:G:H5'	2.04	0.58
25:CY:108:GLU:O	25:CY:111:ARG:N	2.36	0.58
25:CY:125:GLY:O	25:CY:128:ALA:HB3	2.04	0.58
35:DA:1337:G:H2'	35:DA:1338:G:O4'	2.04	0.58
35:DA:1423:G:H2'	35:DA:1424:G:H8	1.69	0.58
27:D1:41:ARG:HH22	35:DA:205:G:H1	1.50	0.58
35:DA:2892:A:C5	35:DA:2893:G:H1'	2.38	0.58
35:DA:535:C:O2'	35:DA:536:A:H5'	2.04	0.58
38:DD:134:ARG:HB2	38:DD:135:PHE:CD1	2.37	0.58
39:DE:104:VAL:O	39:DE:167:VAL:HG12	2.02	0.58
41:DG:12:TYR:H	41:DG:12:TYR:HD1	1.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:12:PRO:O	42:DH:13:LYS:HB2	2.03	0.58
43:DI:94:ALA:C	43:DI:96:ASP:H	2.05	0.58
45:DO:86:ILE:CD1	45:DO:86:ILE:N	2.66	0.58
47:DQ:103:MET:CE	47:DQ:125:LEU:HD21	2.34	0.58
48:DR:9:LYS:NZ	48:DR:39:PRO:HA	2.19	0.58
51:DU:49:HIS:O	51:DU:52:ARG:HB2	2.03	0.58
53:DW:1:MET:HG3	53:DW:2:GLU:N	2.19	0.58
54:DX:12:VAL:HG13	54:DX:17:ALA:CB	2.34	0.58
55:DY:11:ASP:N	55:DY:27:VAL:HG22	2.19	0.58
2:AB:97:TRP:NE1	2:AB:101:MET:SD	2.77	0.57
2:AB:212:GLN:CG	2:AB:235:SER:HB2	2.33	0.57
2:AB:68:ILE:HG22	2:AB:70:PHE:HD1	1.67	0.57
8:AH:96:GLY:O	8:AH:98:LYS:N	2.36	0.57
11:AK:20:TYR:O	11:AK:30:VAL:HA	2.04	0.57
20:AT:26:ASN:HD22	20:AT:26:ASN:N	2.02	0.57
25:AY:150:SER:C	25:AY:152:ASP:H	2.04	0.57
35:BA:1417:C:O2'	35:BA:1418:G:H5'	2.04	0.57
35:BA:1486:A:H61	35:BA:1504:C:H42	1.52	0.57
35:BA:1822:G:H2'	35:BA:1823:G:H8	1.67	0.57
35:BA:2052:G:N3	39:BE:149:ARG:HA	2.19	0.57
35:BA:2307:G:H21	35:BA:2308:G:H5'	1.68	0.57
35:BA:2846:G:H2'	35:BA:2847:U:C6	2.39	0.57
35:BA:319:C:O2'	35:BA:320:A:H5'	2.04	0.57
35:BA:909:A:H1'	47:BQ:10:ARG:HH22	1.69	0.57
35:BA:916:G:C2'	35:BA:917:A:H5''	2.34	0.57
36:BB:42:C:O4'	41:BG:68:PRO:O	2.21	0.57
35:BA:782:A:H2	38:BD:226:MET:HG2	1.64	0.57
38:BD:9:TYR:O	38:BD:10:THR:HG22	2.04	0.57
39:BE:120:TRP:CE3	39:BE:155:LYS:HD3	2.39	0.57
39:BE:120:TRP:NE1	39:BE:155:LYS:HB3	2.18	0.57
39:BE:9:VAL:HG22	39:BE:25:VAL:HB	1.85	0.57
50:BT:28:VAL:N	50:BT:88:ILE:HD13	2.18	0.57
53:BW:29:LEU:O	53:BW:33:ARG:HG3	2.04	0.57
55:BY:47:LYS:HG3	55:BY:60:PHE:CZ	2.39	0.57
1:CA:363:A:C5	12:CL:31:PRO:HD2	2.39	0.57
1:CA:714:G:H2'	1:CA:715:A:C8	2.39	0.57
1:CA:766:A:H2'	1:CA:767:A:O4'	2.04	0.57
1:CA:857:C:H2'	1:CA:858:G:O4'	2.04	0.57
2:CB:169:LYS:HD3	2:CB:169:LYS:O	2.03	0.57
2:CB:237:ALA:O	2:CB:238:LEU:HB3	2.03	0.57
3:CC:120:VAL:HA	3:CC:123:GLN:HE21	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:20:TYR:HE2	8:CH:75:ARG:HD2	1.69	0.57
8:CH:63:LEU:HG	8:CH:65:TYR:OH	2.04	0.57
12:CL:26:ALA:O	12:CL:27:LEU:HB2	2.03	0.57
13:CM:115:LYS:O	13:CM:117:VAL:HG23	2.04	0.57
20:CT:17:ARG:HA	20:CT:20:LEU:HD12	1.86	0.57
20:CT:26:ASN:HD22	20:CT:26:ASN:N	2.00	0.57
26:D0:27:GLU:HG3	26:D0:68:GLU:HA	1.86	0.57
27:D1:23:LYS:HB3	27:D1:23:LYS:HZ2	1.68	0.57
35:DA:1292:U:H2'	35:DA:1293:C:C6	2.39	0.57
35:DA:1642:G:H2'	35:DA:1643:G:C8	2.39	0.57
35:DA:2661:G:H2'	35:DA:2662:A:H8	1.68	0.57
35:DA:40:C:H2'	35:DA:41:C:H6	1.68	0.57
35:DA:464:U:H2'	35:DA:465:G:C8	2.39	0.57
35:DA:590:A:H2'	35:DA:591:C:C6	2.39	0.57
39:DE:12:THR:CG2	50:DT:8:LYS:HE2	2.34	0.57
39:DE:169:ASN:OD1	39:DE:201:THR:HG21	2.03	0.57
39:DE:36:ARG:NH2	39:DE:88:GLY:CA	2.66	0.57
39:DE:59:VAL:HG22	39:DE:63:LEU:HA	1.86	0.57
41:DG:103:LEU:O	41:DG:107:LEU:N	2.37	0.57
41:DG:178:PHE:HB3	41:DG:180:PHE:CE1	2.32	0.57
44:DN:78:TYR:N	44:DN:79:PRO:HD2	2.19	0.57
45:DO:80:ASP:O	45:DO:81:ASP:HB3	2.04	0.57
51:DU:34:LYS:HA	51:DU:34:LYS:CE	2.33	0.57
56:DZ:146:ILE:HA	56:DZ:174:VAL:CG1	2.33	0.57
1:AA:148:G:H2'	1:AA:149:A:C8	2.36	0.57
1:AA:431:A:H2'	1:AA:432:A:C8	2.40	0.57
2:AB:166:ASP:HB2	2:AB:205:ASP:OD2	2.05	0.57
3:AC:153:VAL:HG12	3:AC:154:SER:N	2.19	0.57
4:AD:193:ASP:HB2	4:AD:194:LEU:HD22	1.86	0.57
4:AD:196:LEU:N	4:AD:196:LEU:HD12	2.19	0.57
5:AE:87:SER:OG	5:AE:125:SER:HB3	2.03	0.57
11:AK:61:ALA:CB	11:AK:90:GLY:HA3	2.32	0.57
1:AA:1226:C:H2'	13:AM:103:THR:OG1	2.04	0.57
16:AP:6:LEU:HB3	16:AP:17:TYR:HB3	1.87	0.57
20:AT:42:GLN:O	20:AT:43:LEU:HD23	2.04	0.57
28:B2:22:GLU:HA	28:B2:25:VAL:HG12	1.85	0.57
29:B3:4:LEU:HD12	29:B3:39:ASP:OD1	2.04	0.57
33:B7:30:VAL:HG12	33:B7:33:ARG:HH12	1.69	0.57
35:BA:1014:U:H2'	35:BA:1015:G:H8	1.68	0.57
35:BA:1568:G:H4'	38:BD:59:LYS:HG3	1.86	0.57
35:BA:1658:C:H2'	35:BA:1659:U:C6	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1695:G:H2'	35:BA:1696:G:O4'	2.03	0.57
35:BA:1937:A:O2'	35:BA:1938:A:OP1	2.20	0.57
35:BA:200:U:H2'	35:BA:201:C:H5'	1.86	0.57
35:BA:2090:G:C6	35:BA:2091:U:C4	2.92	0.57
35:BA:2315:G:H2'	35:BA:2316:C:H6	1.67	0.57
35:BA:2843:G:H1	35:BA:2874:C:N4	2.03	0.57
35:BA:380:U:H2'	35:BA:381:G:C8	2.37	0.57
35:BA:913:U:H4'	35:BA:914:C:OP1	2.04	0.57
37:BC:56:GLN:HE22	37:BC:169:GLY:H	1.51	0.57
38:BD:160:GLY:H	38:BD:196:VAL:HB	1.69	0.57
38:BD:218:ARG:HB3	38:BD:219:PRO:HD2	1.84	0.57
41:BG:138:GLN:HG2	41:BG:153:ARG:H	1.68	0.57
41:BG:51:ARG:NE	41:BG:51:ARG:HA	2.19	0.57
41:BG:37:VAL:HB	41:BG:94:LEU:HB2	1.85	0.57
42:BH:89:ILE:N	42:BH:89:ILE:HD13	2.19	0.57
43:BI:6:LEU:O	43:BI:7:GLU:C	2.43	0.57
44:BN:46:VAL:CG1	44:BN:47:ALA:H	2.00	0.57
46:BP:140:ALA:O	46:BP:141:ALA:HB3	2.04	0.57
46:BP:92:GLU:HG3	46:BP:93:GLY:N	2.19	0.57
47:BQ:114:ALA:C	47:BQ:116:GLU:H	2.05	0.57
47:BQ:83:MET:CG	47:BQ:83:MET:O	2.51	0.57
47:BQ:87:LYS:O	47:BQ:88:GLY:O	2.22	0.57
49:BS:61:ASN:ND2	49:BS:62:LYS:HE3	2.18	0.57
51:BU:98:LEU:O	51:BU:101:ARG:N	2.37	0.57
51:BU:92:ARG:HB3	52:BV:11:GLN:CD	2.25	0.57
55:BY:31:LEU:HB2	55:BY:36:ALA:O	2.04	0.57
56:BZ:11:GLU:HB2	56:BZ:13:GLU:HG3	1.85	0.57
1:CA:359:U:H2'	1:CA:360:A:C8	2.39	0.57
4:CD:155:LEU:O	4:CD:158:ILE:HG22	2.03	0.57
12:CL:100:ILE:O	12:CL:102:ARG:N	2.36	0.57
13:CM:82:MET:HB3	13:CM:93:ARG:HH11	1.69	0.57
34:D8:39:LYS:NZ	34:D8:43:GLN:HG3	2.19	0.57
35:DA:1114:G:C2'	35:DA:1115:G:H5''	2.32	0.57
35:DA:1581:G:H2'	35:DA:1582:C:O4'	2.04	0.57
35:DA:2340:G:O2'	35:DA:2341:G:H5'	2.04	0.57
35:DA:2836:U:H2'	35:DA:2837:G:C8	2.38	0.57
35:DA:45:C:H2'	35:DA:47:C:C6	2.39	0.57
35:DA:558:G:H2'	35:DA:559:G:C8	2.36	0.57
35:DA:674:G:O2'	40:DF:74:ARG:HB2	2.04	0.57
35:DA:823:G:H2'	35:DA:824:A:C8	2.39	0.57
35:DA:902:C:H2'	35:DA:903:C:C6	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:173:VAL:HG12	38:DD:185:VAL:O	2.04	0.57
40:DF:148:LEU:HD21	40:DF:191:ARG:HD3	1.85	0.57
43:DI:75:LEU:HD11	43:DI:105:HIS:HE1	1.68	0.57
43:DI:71:ILE:HG13	43:DI:72:LEU:H	1.67	0.57
43:DI:83:ALA:O	43:DI:144:VAL:HG13	2.04	0.57
43:DI:94:ALA:HA	43:DI:97:ILE:CG1	2.34	0.57
44:DN:15:LEU:HB3	44:DN:136:GLU:HA	1.84	0.57
46:DP:111:ARG:HG3	46:DP:128:HIS:ND1	2.18	0.57
34:D8:59:LYS:HG3	46:DP:49:ARG:HD2	1.86	0.57
46:DP:90:ARG:CD	46:DP:91:PHE:HD1	2.17	0.57
49:DS:49:VAL:HG21	49:DS:77:ALA:HB2	1.86	0.57
50:DT:83:ILE:HD11	50:DT:84:GLN:HE21	1.67	0.57
51:DU:55:ARG:HA	51:DU:58:ARG:CD	2.34	0.57
54:DX:21:PHE:HE1	54:DX:26:TYR:HB3	1.69	0.57
56:DZ:45:ASP:O	56:DZ:49:ARG:HG2	2.03	0.57
1:AA:1160:G:OP1	2:AB:132:LYS:HE3	2.04	0.57
1:AA:159:G:N1	1:AA:163:C:N4	2.53	0.57
1:AA:599:C:H2'	1:AA:600:C:H6	1.67	0.57
2:AB:102:LEU:H	2:AB:102:LEU:HD12	1.67	0.57
3:AC:120:VAL:HA	3:AC:123:GLN:HE21	1.69	0.57
6:AF:43:LEU:H	6:AF:43:LEU:CD1	2.17	0.57
7:AG:69:VAL:HG22	7:AG:135:VAL:HG22	1.86	0.57
8:AH:91:ARG:HG2	8:AH:91:ARG:HH11	1.69	0.57
10:AJ:94:VAL:HG12	10:AJ:95:GLU:N	2.19	0.57
13:AM:76:ALA:HA	13:AM:79:LYS:HD2	1.85	0.57
16:AP:21:VAL:O	16:AP:33:ILE:HB	2.04	0.57
16:AP:45:THR:HG22	16:AP:47:ASP:N	2.19	0.57
16:AP:64:ALA:O	16:AP:65:GLN:C	2.42	0.57
17:AQ:56:VAL:HG23	17:AQ:78:GLU:O	2.05	0.57
27:B1:87:PRO:CG	27:B1:88:LYS:H	2.16	0.57
34:B8:60:LEU:HA	34:B8:63:PRO:HG2	1.85	0.57
35:BA:1600:C:C2'	35:BA:1601:G:H5'	2.34	0.57
1:AA:1494:G:H5''	35:BA:1913:A:C6	2.40	0.57
35:BA:2040:C:H2'	35:BA:2041:U:C6	2.40	0.57
35:BA:2645:G:C3'	35:BA:2646:C:H5'	2.28	0.57
35:BA:2808:U:H2'	35:BA:2809:A:C5'	2.33	0.57
40:BF:67:GLN:O	40:BF:68:LYS:HG2	2.05	0.57
45:BO:13:ASN:HD22	45:BO:97:ARG:CB	2.18	0.57
46:BP:86:LYS:HB3	46:BP:117:GLU:C	2.24	0.57
48:BR:42:LYS:HG3	48:BR:45:ARG:NH2	2.18	0.57
48:BR:52:ILE:O	48:BR:55:ALA:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:BU:107:ALA:O	51:BU:111:GLU:HG2	2.04	0.57
51:BU:91:ASP:OD2	51:BU:96:ALA:CA	2.53	0.57
55:BY:74:PRO:O	55:BY:75:ILE:CB	2.52	0.57
1:CA:103:C:H2'	1:CA:104:G:H8	1.70	0.57
1:CA:1248:A:H2'	1:CA:1249:C:H5'	1.85	0.57
1:CA:1406:U:O2'	1:CA:1407:C:H5'	2.04	0.57
1:CA:877:C:H5''	8:CH:88:LYS:HD2	1.86	0.57
4:CD:176:LEU:CG	4:CD:177:ASP:N	2.66	0.57
4:CD:180:GLY:C	4:CD:181:MET:HG2	2.24	0.57
1:CA:18:C:P	5:CE:127:ASN:HD21	2.27	0.57
6:CF:27:GLN:HE21	6:CF:27:GLN:HA	1.69	0.57
7:CG:95:ARG:O	7:CG:96:GLN:C	2.43	0.57
1:CA:600:C:H5'	8:CH:129:VAL:O	2.05	0.57
11:CK:86:GLY:H	11:CK:112:THR:CG2	2.17	0.57
12:CL:90:VAL:O	12:CL:92:ASP:N	2.34	0.57
14:CN:51:GLY:O	14:CN:53:LEU:N	2.36	0.57
16:CP:14:ASN:N	16:CP:15:PRO:CD	2.67	0.57
16:CP:71:ARG:NH1	16:CP:71:ARG:HG3	2.20	0.57
17:CQ:12:SER:HB3	17:CQ:20:THR:OG1	2.04	0.57
18:CR:86:VAL:HG12	18:CR:87:ARG:NH1	2.19	0.57
13:CM:87:TYR:N	19:CS:73:GLU:O	2.33	0.57
28:D2:12:GLU:CD	28:D2:14:ARG:NH1	2.58	0.57
32:D6:14:THR:O	32:D6:49:HIS:HA	2.04	0.57
33:D7:34:ARG:O	33:D7:35:ARG:C	2.41	0.57
35:DA:812:C:C2	35:DA:1250:G:N1	2.72	0.57
35:DA:1298:C:H2'	35:DA:1299:G:C8	2.39	0.57
35:DA:139:G:H1	35:DA:142(A):C:H42	1.51	0.57
35:DA:1999:C:H4'	35:DA:2723:C:O2	2.05	0.57
35:DA:2364:C:O2'	35:DA:2365:G:H5'	2.03	0.57
35:DA:692:C:H2'	35:DA:693:C:H6	1.69	0.57
35:DA:763:G:C4	35:DA:765:G:C8	2.93	0.57
40:DF:31:HIS:O	40:DF:32:LEU:C	2.43	0.57
41:DG:49:ASP:O	41:DG:50:ALA:HB2	2.04	0.57
41:DG:47:LYS:O	41:DG:51:ARG:HG2	2.04	0.57
42:DH:102:ALA:CB	42:DH:117:PRO:HD3	2.21	0.57
42:DH:20:ALA:HB1	42:DH:21:PRO:CD	2.34	0.57
42:DH:19:VAL:CG2	42:DH:44:VAL:HA	2.32	0.57
42:DH:89:ILE:CD1	42:DH:90:LYS:H	2.17	0.57
44:DN:66:LYS:O	44:DN:87:LEU:HB3	2.05	0.57
46:DP:39:LYS:HD3	46:DP:40:SER:N	2.19	0.57
46:DP:92:GLU:HG3	46:DP:93:GLY:N	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:141:GLN:NE2	56:DZ:89:PHE:HB3	2.19	0.57
47:DQ:68:ILE:HD13	47:DQ:68:ILE:N	2.15	0.57
50:DT:89:VAL:HG11	50:DT:91:ARG:HE	1.68	0.57
51:DU:31:SER:O	51:DU:33:ARG:N	2.37	0.57
52:DV:40:LEU:O	52:DV:40:LEU:HD13	2.05	0.57
1:AA:1508:G:H2'	1:AA:1509:C:H6	1.70	0.57
1:AA:341:C:O2'	1:AA:342:C:H5'	2.04	0.57
1:AA:404:U:H2'	1:AA:405:U:H6	1.68	0.57
1:AA:569:C:H42	1:AA:881:G:H1	1.53	0.57
2:AB:71:VAL:CG2	2:AB:93:VAL:HG23	2.32	0.57
4:AD:154:ASN:CB	4:AD:159:ARG:HH21	2.17	0.57
5:AE:15:ARG:HG3	5:AE:28:PHE:CE2	2.39	0.57
8:AH:119:LEU:HD12	8:AH:124:ALA:N	2.18	0.57
10:AJ:48:THR:HG23	10:AJ:61:GLU:C	2.25	0.57
19:AS:41:VAL:HB	19:AS:44:MET:CB	2.35	0.57
20:AT:19:SER:O	20:AT:23:ARG:N	2.35	0.57
20:AT:32:ALA:O	20:AT:36:LEU:HD23	2.04	0.57
25:AY:60:ALA:HB2	25:AY:66:LEU:HG	1.86	0.57
27:B1:73:LEU:O	27:B1:76:ARG:HG2	2.03	0.57
28:B2:56:GLN:OE1	35:BA:76:C:H1'	2.04	0.57
35:BA:1821:A:H2'	35:BA:1822:G:H8	1.68	0.57
35:BA:1910:G:C6	35:BA:1921:G:C6	2.93	0.57
35:BA:528:A:C2	35:BA:2043:C:C5'	2.83	0.57
35:BA:2469:A:H3'	35:BA:2470:G:O4'	2.04	0.57
35:BA:588:U:O4	35:BA:670:A:O2'	2.21	0.57
35:BA:827:U:H2'	35:BA:2068:U:N3	2.20	0.57
38:BD:183:ARG:HG2	38:BD:183:ARG:NH1	2.18	0.57
38:BD:34:VAL:HG22	38:BD:35:LYS:HZ2	1.69	0.57
41:BG:85:GLY:O	41:BG:87:PRO:HD3	2.05	0.57
46:BP:112:LEU:O	46:BP:128:HIS:HB2	2.04	0.57
50:BT:86:ILE:HG12	50:BT:87:ASP:O	2.04	0.57
51:BU:102:GLU:O	51:BU:105:VAL:HG23	2.04	0.57
52:BV:19:LYS:HZ2	52:BV:20:LEU:N	1.91	0.57
52:BV:4:ILE:HD12	52:BV:40:LEU:HD11	1.86	0.57
35:BA:1224:C:O3'	52:BV:88:ARG:HB3	2.05	0.57
1:CA:1057:G:H5''	3:CC:154:SER:OG	2.04	0.57
1:CA:1298:C:H1'	1:CA:1299:A:C2	2.39	0.57
1:CA:815:A:N7	1:CA:1509:C:O2'	2.37	0.57
1:CA:1511:G:H8	1:CA:1511:G:O5'	1.86	0.57
1:CA:409:G:H2'	1:CA:410:G:O4'	2.04	0.57
1:CA:625:G:C4	1:CA:626:U:C5	2.92	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:83:U:H2'	1:CA:84:U:C6	2.40	0.57
1:CA:899:C:H2'	1:CA:900:A:O4'	2.04	0.57
2:CB:187:LEU:CA	2:CB:201:ILE:HB	2.34	0.57
2:CB:39:ILE:CG2	2:CB:40:HIS:N	2.66	0.57
4:CD:118:ARG:O	4:CD:121:VAL:HB	2.04	0.57
5:CE:103:GLY:O	5:CE:106:PRO:HD2	2.04	0.57
1:CA:688:G:H5'	11:CK:46:GLY:O	2.03	0.57
11:CK:21:ILE:CB	11:CK:84:VAL:HG12	2.33	0.57
13:CM:77:ASN:O	13:CM:81:LEU:HG	2.05	0.57
18:CR:87:ARG:CZ	18:CR:87:ARG:HB3	2.34	0.57
25:CY:126:ARG:O	25:CY:127:VAL:C	2.43	0.57
25:CY:6:LEU:O	25:CY:6:LEU:HD22	2.05	0.57
26:D0:27:GLU:HB2	26:D0:69:PHE:HD1	1.66	0.57
27:D1:66:HIS:C	27:D1:68:PRO:HD2	2.24	0.57
34:D8:38:GLY:C	34:D8:40:GLU:H	2.07	0.57
35:DA:1418:G:N1	35:DA:1579:A:H5'	2.19	0.57
35:DA:45:C:OP2	35:DA:215:G:H2'	2.05	0.57
35:DA:2455:G:H2'	35:DA:2456:C:C6	2.39	0.57
35:DA:540:C:H2'	35:DA:541:C:C5	2.39	0.57
35:DA:836:G:H2'	35:DA:837:C:H6	1.67	0.57
39:DE:79:ARG:HH11	39:DE:79:ARG:HG2	1.68	0.57
42:DH:109:PHE:CE1	42:DH:152:ARG:NE	2.72	0.57
44:DN:78:TYR:CD1	44:DN:79:PRO:HD3	2.39	0.57
45:DO:88:ASN:OD1	45:DO:92:GLU:N	2.31	0.57
46:DP:50:ARG:CZ	46:DP:51:PHE:CZ	2.88	0.57
49:DS:83:LYS:HE3	49:DS:84:GLN:HG3	1.87	0.57
50:DT:102:ILE:CA	50:DT:110:ILE:HD11	2.34	0.57
50:DT:23:ARG:NH2	50:DT:120:ARG:HD3	2.20	0.57
51:DU:92:ARG:CG	51:DU:94:ASN:HB3	2.35	0.57
52:DV:38:LEU:CD2	52:DV:39:LEU:N	2.63	0.57
54:DX:36:LYS:HD3	54:DX:38:GLU:HB2	1.86	0.57
55:DY:16:ALA:HA	55:DY:21:LYS:HD2	1.86	0.57
1:AA:1175:G:H2'	1:AA:1176:A:C8	2.40	0.57
1:AA:1190:G:C8	3:AC:3:ASN:ND2	2.71	0.57
1:AA:409:G:H2'	1:AA:410:G:O4'	2.03	0.57
1:AA:634:C:O2'	1:AA:635:G:H5'	2.03	0.57
2:AB:39:ILE:CG2	2:AB:40:HIS:N	2.67	0.57
4:AD:153:ARG:HB3	4:AD:153:ARG:HH11	1.69	0.57
4:AD:176:LEU:HD21	4:AD:178:VAL:HG22	1.84	0.57
7:AG:143:ARG:CB	7:AG:143:ARG:HH11	2.18	0.57
1:AA:972:C:H5'	10:AJ:57:LYS:NZ	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:99:GLN:HE22	11:AK:105:VAL:HG11	1.69	0.57
16:AP:60:LEU:C	16:AP:62:VAL:H	2.08	0.57
25:AY:17:SER:HB3	25:AY:132:ILE:HD11	1.87	0.57
27:B1:76:ARG:HA	27:B1:76:ARG:NE	2.18	0.57
28:B2:32:LEU:HD11	28:B2:33:MET:SD	2.45	0.57
29:B3:31:LEU:CD2	29:B3:32:GLN:H	2.16	0.57
34:B8:11:LYS:HG2	34:B8:11:LYS:O	2.04	0.57
34:B8:51:ALA:HA	34:B8:54:GLU:OE1	2.03	0.57
35:BA:1108:U:C2'	35:BA:1109:C:H5'	2.34	0.57
35:BA:2019:A:O3'	51:BU:27:LEU:HD12	2.05	0.57
35:BA:2579:C:C2'	35:BA:2580:U:H5'	2.34	0.57
35:BA:2683:C:H5''	50:BT:53:ARG:HH22	1.67	0.57
35:BA:2802:G:O2'	35:BA:2803:C:C5'	2.52	0.57
35:BA:30:G:O2'	35:BA:31:C:H5'	2.05	0.57
35:BA:618:C:H2'	35:BA:619:G:O4'	2.04	0.57
35:BA:692:C:H2'	35:BA:693:C:C6	2.40	0.57
39:BE:200:GLU:OE2	39:BE:200:GLU:N	2.37	0.57
39:BE:36:ARG:HA	39:BE:46:ALA:O	2.04	0.57
40:BF:167:ALA:HB1	40:BF:173:VAL:HG11	1.87	0.57
40:BF:24:LEU:HB3	40:BF:25:PRO:CD	2.35	0.57
43:BI:75:LEU:HD11	43:BI:105:HIS:HE1	1.70	0.57
44:BN:74:ARG:NH2	44:BN:101:HIS:O	2.35	0.57
46:BP:16:ARG:CD	46:BP:16:ARG:C	2.73	0.57
49:BS:25:ARG:HH21	49:BS:89:ARG:NH1	1.96	0.57
49:BS:26:LEU:HD22	49:BS:87:PHE:CE1	2.39	0.57
53:BW:37:ARG:HG3	53:BW:37:ARG:HH11	1.69	0.57
54:BX:55:ASN:O	54:BX:77:LYS:CB	2.53	0.57
2:CB:55:PHE:CD1	2:CB:58:ILE:HD12	2.40	0.57
1:CA:1056:U:H5'	3:CC:163:ALA:CB	2.34	0.57
6:CF:30:LEU:HB3	6:CF:35:ALA:CB	2.34	0.57
6:CF:76:ALA:HB1	6:CF:80:ARG:HH21	1.68	0.57
11:CK:20:TYR:O	11:CK:30:VAL:HA	2.05	0.57
12:CL:37:CYS:SG	12:CL:81:SER:HB2	2.45	0.57
17:CQ:58:GLU:HB2	17:CQ:74:LEU:HB3	1.86	0.57
26:D0:53:MET:HG3	26:D0:59:LEU:HD23	1.86	0.57
28:D2:14:ARG:O	28:D2:16:LEU:N	2.38	0.57
31:D5:43:HIS:CD2	31:D5:43:HIS:N	2.71	0.57
35:DA:1451:C:H4'	35:DA:1452:A:C8	2.39	0.57
35:DA:1695:G:H2'	35:DA:1696:G:O4'	2.05	0.57
35:DA:2881:C:C2	35:DA:2882:A:C8	2.92	0.57
35:DA:292:C:N4	35:DA:348:G:H1	2.01	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:455:C:H3'	35:DA:456:C:H5''	1.86	0.57
35:DA:514:A:H1'	35:DA:581:C:O2'	2.04	0.57
38:DD:169:GLU:O	38:DD:171:ASP:N	2.36	0.57
40:DF:65:TRP:HH2	40:DF:75:HIS:HD2	1.49	0.57
43:DI:58:LEU:C	43:DI:60:GLU:H	2.08	0.57
44:DN:72:TYR:HB3	44:DN:74:ARG:HG2	1.86	0.57
35:DA:587:C:C4	46:DP:33:ARG:HD2	2.39	0.57
48:DR:60:LEU:O	48:DR:61:HIS:C	2.43	0.57
52:DV:52:VAL:O	52:DV:54:GLY:N	2.37	0.57
53:DW:1:MET:HA	53:DW:1:MET:HE3	1.84	0.57
54:DX:32:PRO:HD3	54:DX:72:LYS:NZ	2.19	0.57
56:DZ:144:LEU:HD11	56:DZ:150:LEU:HD13	1.87	0.57
1:AA:178:C:H2'	1:AA:179:A:H8	1.69	0.57
2:AB:137:ARG:HG2	2:AB:137:ARG:NH1	2.20	0.57
2:AB:187:LEU:CA	2:AB:201:ILE:HB	2.34	0.57
4:AD:22:LYS:HB2	4:AD:26:CYS:CB	2.24	0.57
7:AG:84:ASN:ND2	7:AG:84:ASN:N	2.53	0.57
15:AO:6:GLU:CD	15:AO:6:GLU:H	2.08	0.57
21:AU:17:THR:O	21:AU:22:ARG:HD3	2.05	0.57
23:AW:39:A:C2'	23:AW:40:C:H5'	2.35	0.57
23:AW:51:U:H3	23:AW:65:G:H1	1.51	0.57
25:AY:28:LEU:O	25:AY:30:THR:HG23	2.03	0.57
25:AY:29:ARG:HB3	25:AY:32:ARG:NH2	2.18	0.57
25:AY:75:ALA:O	25:AY:79:ILE:HG13	2.03	0.57
27:B1:48:LYS:HG3	27:B1:49:VAL:N	2.13	0.57
28:B2:43:GLN:O	28:B2:46:GLN:HB3	2.04	0.57
35:BA:1009:A:H2'	35:BA:1010:A:C8	2.39	0.57
35:BA:1197:G:H2'	35:BA:1198:U:H6	1.69	0.57
35:BA:1292:U:H2'	35:BA:1293:C:C6	2.40	0.57
35:BA:2081:C:H2'	35:BA:2082:A:H8	1.68	0.57
35:BA:2380:C:H2'	35:BA:2381:C:C6	2.40	0.57
35:BA:2600:A:C2'	35:BA:2601:C:H5'	2.33	0.57
35:BA:2666:C:H2'	35:BA:2667:C:O4'	2.05	0.57
35:BA:2839:G:H21	48:BR:92:GLY:HA3	1.70	0.57
37:BC:68:LEU:HB3	37:BC:70:LYS:HG2	1.86	0.57
38:BD:134:ARG:HB2	38:BD:135:PHE:CD1	2.40	0.57
35:BA:2619:C:OP1	39:BE:152:LYS:HD3	2.04	0.57
41:BG:173:LEU:HD12	41:BG:178:PHE:CZ	2.40	0.57
42:BH:138:LYS:H	42:BH:141:VAL:HG23	1.69	0.57
46:BP:50:ARG:HD2	46:BP:51:PHE:CG	2.39	0.57
47:BQ:20:ALA:CA	47:BQ:98:LYS:HB3	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:78:LEU:O	50:BT:79:HIS:ND1	2.37	0.57
50:BT:89:VAL:CG1	50:BT:91:ARG:HE	2.17	0.57
51:BU:106:PHE:O	51:BU:110:VAL:HG23	2.04	0.57
56:BZ:7:ALA:O	56:BZ:38:TYR:O	2.22	0.57
1:CA:1222:G:H5''	19:CS:78:ARG:HH11	1.69	0.57
1:CA:1223:C:P	1:CA:1224:G:H2'	2.44	0.57
1:CA:1524:C:H5''	11:CK:120:ARG:HH12	1.69	0.57
1:CA:176:C:H2'	1:CA:177:C:H6	1.68	0.57
1:CA:272:C:O2'	1:CA:273:A:H5'	2.04	0.57
1:CA:644:G:O2'	1:CA:645:C:H5'	2.04	0.57
1:CA:708:C:O2'	1:CA:709:G:H5'	2.05	0.57
2:CB:171:ALA:O	2:CB:174:VAL:HB	2.05	0.57
4:CD:133:VAL:HG12	4:CD:135:LEU:H	1.69	0.57
4:CD:193:ASP:HB2	4:CD:194:LEU:HD22	1.85	0.57
6:CF:78:GLU:O	6:CF:81:ILE:HG13	2.05	0.57
9:CI:5:TYR:HA	9:CI:17:VAL:O	2.05	0.57
19:CS:49:ILE:O	19:CS:60:VAL:HG12	2.04	0.57
20:CT:16:HIS:O	20:CT:19:SER:HB2	2.04	0.57
20:CT:48:LYS:O	20:CT:49:ALA:HB2	2.04	0.57
25:CY:19:GLU:O	25:CY:22:GLU:HB2	2.05	0.57
27:D1:89:GLU:CD	27:D1:89:GLU:N	2.57	0.57
35:DA:1623:G:H2'	35:DA:1624:G:H8	1.70	0.57
35:DA:1779:U:C2	35:DA:1783:A:N7	2.72	0.57
35:DA:2625:G:H2'	35:DA:2626:C:C6	2.38	0.57
35:DA:2753:A:H2	35:DA:2754:U:C2	2.22	0.57
35:DA:2828:C:H2'	35:DA:2829:C:C6	2.39	0.57
35:DA:325:G:H2'	35:DA:326:G:H8	1.68	0.57
35:DA:464:U:H2'	35:DA:465:G:O4'	2.03	0.57
35:DA:549:G:H2'	35:DA:551:G:O4'	2.05	0.57
38:DD:108:PRO:HG2	38:DD:111:LEU:HB2	1.86	0.57
39:DE:52:LEU:N	39:DE:74:PRO:HB2	2.19	0.57
42:DH:149:ARG:HA	42:DH:162:ILE:CD1	2.34	0.57
35:DA:2748:A:H2	42:DH:63:SER:HB3	1.67	0.57
44:DN:26:LEU:HD21	44:DN:30:ILE:HD11	1.84	0.57
46:DP:16:ARG:CD	46:DP:16:ARG:C	2.73	0.57
46:DP:91:PHE:N	46:DP:91:PHE:CD1	2.72	0.57
47:DQ:87:LYS:O	47:DQ:88:GLY:O	2.22	0.57
52:DV:4:ILE:HD12	52:DV:40:LEU:HD11	1.86	0.57
53:DW:5:ALA:HB3	53:DW:105:VAL:H	1.68	0.57
56:DZ:119:GLU:OE1	56:DZ:122:ARG:HB3	2.04	0.57
1:AA:924:C:H2'	1:AA:925:G:C8	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:959:A:H2'	1:AA:960:U:H4'	1.87	0.57
1:AA:974:A:C1'	14:AN:31:ARG:HH21	2.17	0.57
3:AC:127:ARG:HG2	3:AC:127:ARG:NH1	2.20	0.57
7:AG:36:LYS:HB2	7:AG:36:LYS:HZ2	1.69	0.57
10:AJ:4:ILE:HB	10:AJ:74:ILE:CG1	2.34	0.57
12:AL:119:LYS:O	12:AL:120:TYR:HB2	2.05	0.57
25:AY:21:LEU:O	25:AY:25:LEU:HG	2.05	0.57
26:B0:27:GLU:HB2	26:B0:69:PHE:HD1	1.70	0.57
27:B1:47:GLN:HE21	27:B1:64:ALA:CB	2.18	0.57
32:B6:39:TYR:CE1	35:BA:2347:C:H4'	2.35	0.57
33:B7:30:VAL:HG23	33:B7:31:LEU:N	2.20	0.57
35:BA:1234:U:H2'	35:BA:1234:U:O2	2.04	0.57
35:BA:1399:C:H2'	35:BA:1400:G:H8	1.68	0.57
35:BA:141:A:H8	35:BA:1408:C:HO2'	1.50	0.57
35:BA:1528(A):A:H2'	35:BA:1529:G:H5''	1.85	0.57
35:BA:1820:U:C2	38:BD:202:LYS:HB3	2.40	0.57
35:BA:1999:C:H2'	35:BA:2000:G:C8	2.38	0.57
35:BA:2744:G:O2'	35:BA:2745:C:H5'	2.05	0.57
35:BA:709:U:H2'	35:BA:710:G:C8	2.37	0.57
36:BB:55:U:O2'	36:BB:56:G:H5'	2.05	0.57
37:BC:55:ASP:CG	37:BC:56:GLN:H	2.08	0.57
38:BD:125:ILE:O	38:BD:125:ILE:HG22	2.04	0.57
35:BA:773:U:H4'	38:BD:47:GLY:HA2	1.87	0.57
38:BD:81:ALA:N	38:BD:94:LEU:HD11	2.20	0.57
40:BF:22:ALA:CA	40:BF:26:ALA:HB2	2.35	0.57
40:BF:88:VAL:CG2	40:BF:89:VAL:N	2.67	0.57
41:BG:138:GLN:CG	41:BG:153:ARG:H	2.17	0.57
41:BG:174:GLU:HG3	41:BG:182:LYS:NZ	2.19	0.57
42:BH:159:GLU:OE1	42:BH:159:GLU:HA	2.05	0.57
43:BI:102:SER:HA	43:BI:107:VAL:O	2.04	0.57
43:BI:10:GLU:O	43:BI:11:ASN:HB3	2.04	0.57
46:BP:99:LEU:O	46:BP:102:ARG:HB3	2.04	0.57
46:BP:39:LYS:CD	46:BP:40:SER:N	2.61	0.57
46:BP:70:GLN:CG	46:BP:71:VAL:H	2.16	0.57
47:BQ:32:TYR:HD1	47:BQ:32:TYR:H	1.51	0.57
54:BX:35:THR:O	54:BX:39:ILE:HG23	2.05	0.57
55:BY:68:HIS:HB3	55:BY:71:LYS:CE	2.34	0.57
56:BZ:48:PHE:CE2	56:BZ:52:SER:HA	2.39	0.57
56:BZ:81:ARG:O	56:BZ:82:ARG:C	2.42	0.57
3:CC:138:VAL:CG2	3:CC:151:VAL:HG23	2.35	0.57
4:CD:33:MET:HA	4:CD:33:MET:CE	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CF:87:ARG:HG3	6:CF:87:ARG:HH11	1.69	0.57
9:CI:26:VAL:HG13	9:CI:63:ILE:HD11	1.86	0.57
1:CA:280:C:O2	17:CQ:38:ARG:HG3	2.04	0.57
28:D2:12:GLU:OE1	28:D2:12:GLU:O	2.23	0.57
35:DA:1006:C:H2'	35:DA:1007:C:C6	2.40	0.57
35:DA:1523:U:H2'	35:DA:1524:G:C8	2.39	0.57
35:DA:1819:A:H1'	35:DA:1821:A:C6	2.40	0.57
35:DA:2589:A:H2'	35:DA:2590:A:H8	1.69	0.57
35:DA:619:G:P	35:DA:620:G:H22	2.26	0.57
35:DA:962:G:O2'	35:DA:963:U:H5'	2.04	0.57
38:DD:204:ILE:O	38:DD:204:ILE:HG13	2.05	0.57
38:DD:259:THR:O	38:DD:260:ARG:C	2.42	0.57
35:DA:2512:C:H4'	39:DE:122:PHE:CZ	2.40	0.57
40:DF:41:LEU:HD23	40:DF:44:ARG:HD3	1.87	0.57
45:DO:22:ILE:H	45:DO:41:ALA:HA	1.70	0.57
46:DP:112:LEU:O	46:DP:128:HIS:HB2	2.05	0.57
46:DP:148:LEU:H	46:DP:148:LEU:HD13	1.70	0.57
46:DP:83:VAL:HG23	46:DP:105:LEU:HD22	1.87	0.57
48:DR:28:LEU:C	48:DR:30:THR:H	2.08	0.57
50:DT:85:LYS:HG2	50:DT:85:LYS:O	2.05	0.57
54:DX:83:VAL:O	54:DX:85:PRO:HD3	2.05	0.57
1:AA:1388:C:H2'	1:AA:1389:C:H6	1.69	0.57
1:AA:1491:G:H5''	1:AA:1492:A:OP1	2.04	0.57
1:AA:1507:A:N1	1:AA:1530:G:C4	2.72	0.57
1:AA:637:G:H2'	1:AA:638:G:H8	1.69	0.57
4:AD:8:VAL:O	4:AD:10:ARG:HB3	2.05	0.57
6:AF:68:PRO:CG	6:AF:71:ARG:HE	2.18	0.57
7:AG:42:ILE:HG23	7:AG:117:ALA:HB2	1.86	0.57
7:AG:36:LYS:NZ	7:AG:36:LYS:HB2	2.19	0.57
8:AH:97:VAL:HG13	8:AH:98:LYS:N	2.14	0.57
13:AM:58:GLU:OE1	13:AM:58:GLU:HA	2.04	0.57
10:AJ:62:HIS:O	14:AN:59:ALA:HB3	2.04	0.57
15:AO:70:LEU:HD23	15:AO:78:TYR:HA	1.87	0.57
23:AW:6:G:H1	23:AW:68:C:H42	1.52	0.57
26:B0:43:THR:HB	26:B0:57:PHE:CE1	2.40	0.57
27:B1:13:ILE:HG23	27:B1:14:VAL:H	1.70	0.57
32:B6:30:THR:HG21	35:BA:2286:A:OP1	2.04	0.57
35:BA:149:A:H2'	35:BA:150:C:O4'	2.05	0.57
35:BA:845:G:O2'	35:BA:846:C:H5	1.88	0.57
35:BA:996:A:H2'	35:BA:997:G:H8	1.69	0.57
38:BD:19:ALA:O	38:BD:21:PHE:CE1	2.58	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:70:TRP:CH2	38:BD:150:LYS:HA	2.39	0.57
42:BH:96:ALA:HB2	42:BH:105:LEU:HB3	1.86	0.57
42:BH:38:SER:O	42:BH:40:GLU:N	2.38	0.57
45:BO:89:ASN:C	45:BO:91:LEU:H	2.07	0.57
50:BT:102:ILE:CA	50:BT:110:ILE:HD11	2.35	0.57
51:BU:106:PHE:CA	51:BU:109:LEU:HD12	2.30	0.57
35:BA:445:C:O3'	51:BU:3:ARG:HG2	2.05	0.57
51:BU:92:ARG:HH22	52:BV:10:LYS:HA	1.70	0.57
28:B2:23:LYS:CA	54:BX:5:TYR:HE1	2.17	0.57
56:BZ:10:ARG:NE	56:BZ:36:LYS:HB3	2.19	0.57
1:CA:1014:A:H2'	1:CA:1015:A:C8	2.40	0.57
1:CA:1190:G:C8	3:CC:3:ASN:ND2	2.70	0.57
1:CA:1268:A:H4'	21:CU:20:LYS:N	2.19	0.57
1:CA:1444:C:N4	1:CA:1458:G:H1	2.03	0.57
1:CA:582:U:H2'	1:CA:583:A:C8	2.40	0.57
2:CB:97:TRP:NE1	2:CB:101:MET:SD	2.78	0.57
2:CB:71:VAL:O	2:CB:164:VAL:HA	2.05	0.57
4:CD:170:VAL:HG13	4:CD:171:GLY:N	2.18	0.57
6:CF:45:LEU:O	6:CF:46:ARG:HD2	2.04	0.57
18:CR:58:LEU:CD1	18:CR:58:LEU:H	2.10	0.57
23:CW:29:C:O2'	23:CW:30:G:H5'	2.05	0.57
25:CY:17:SER:O	25:CY:20:VAL:HB	2.05	0.57
31:D5:20:ARG:HH12	53:DW:15:ARG:HH21	1.53	0.57
31:D5:52:TYR:HA	31:D5:56:LYS:HZ2	1.69	0.57
35:DA:1275:A:C4	48:DR:16:HIS:CE1	2.92	0.57
35:DA:1291:C:O2'	35:DA:1292:U:H5'	2.03	0.57
35:DA:1763:G:OP1	35:DA:1763:G:H4'	2.04	0.57
35:DA:1902:C:C5'	38:DD:246:PRO:HD3	2.34	0.57
35:DA:1993:U:O2'	35:DA:1994:C:H5'	2.05	0.57
35:DA:1998:G:H2'	35:DA:1999:C:H6	1.70	0.57
35:DA:2182:G:H2'	35:DA:2183:C:C6	2.40	0.57
35:DA:2685:G:C2	35:DA:2686:G:N7	2.73	0.57
35:DA:793:A:OP2	35:DA:2071:A:O2'	2.20	0.57
40:DF:20:LEU:HB3	40:DF:23:ASP:OD2	2.04	0.57
41:DG:38:VAL:HB	41:DG:158:ALA:CB	2.35	0.57
45:DO:79:PHE:CE2	45:DO:101:PRO:HB2	2.31	0.57
50:DT:58:ASN:N	50:DT:58:ASN:HD22	2.01	0.57
35:DA:559:G:H22	51:DU:49:HIS:CD2	2.23	0.57
51:DU:55:ARG:HA	51:DU:58:ARG:HD2	1.87	0.57
52:DV:38:LEU:O	52:DV:39:LEU:HD22	2.05	0.57
54:DX:12:VAL:CG1	54:DX:27:THR:HG23	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:68:HIS:ND1	55:DY:69:ALA:N	2.52	0.57
56:DZ:104:PHE:CD1	56:DZ:139:VAL:HG11	2.40	0.57
56:DZ:23:LYS:HD2	56:DZ:38:TYR:HD1	1.70	0.57
56:DZ:10:ARG:HG3	56:DZ:38:TYR:N	2.20	0.57
1:AA:227:G:H2'	1:AA:228:A:C8	2.40	0.57
1:AA:355:C:C4	1:AA:356:A:N7	2.73	0.57
1:AA:511:C:HO2'	1:AA:512:U:H6	1.52	0.57
1:AA:83:U:H2'	1:AA:84:U:C6	2.40	0.57
1:AA:927:G:H2'	1:AA:928:G:H8	1.70	0.57
2:AB:121:LEU:O	2:AB:121:LEU:HD23	2.04	0.57
2:AB:171:ALA:O	2:AB:174:VAL:HB	2.05	0.57
3:AC:113:ALA:HB3	3:AC:114:PRO:HD3	1.86	0.57
5:AE:80:ILE:HG13	5:AE:91:LEU:HB2	1.87	0.57
9:AI:26:VAL:HG13	9:AI:63:ILE:CD1	2.35	0.57
10:AJ:6:ILE:O	10:AJ:6:ILE:HD12	2.04	0.57
11:AK:28:THR:HG22	11:AK:29:ILE:N	2.20	0.57
11:AK:96:ARG:CA	11:AK:99:GLN:HG2	2.33	0.57
20:AT:76:ALA:O	20:AT:80:ARG:HG2	2.05	0.57
27:B1:76:ARG:O	27:B1:77:ALA:CB	2.52	0.57
35:BA:1114:G:C2'	35:BA:1115:G:H5''	2.33	0.57
35:BA:1385:G:H4'	35:BA:1386:C:OP1	2.05	0.57
35:BA:1884:A:H2'	35:BA:1885:A:C5'	2.21	0.57
35:BA:2574:G:H2'	35:BA:2575:C:H6	1.70	0.57
35:BA:2732:G:H3'	35:BA:2733:A:H5'	1.86	0.57
35:BA:2809:A:O2'	35:BA:2810:A:H5'	2.04	0.57
35:BA:455:C:H3'	35:BA:456:C:H5''	1.87	0.57
35:BA:568:U:H2'	35:BA:570:G:OP2	2.05	0.57
35:BA:61:G:O2'	35:BA:62:C:H5'	2.04	0.57
35:BA:695:G:N2	35:BA:696:G:H1'	2.19	0.57
35:BA:782:A:C2	38:BD:226:MET:CG	2.80	0.57
35:BA:809:G:O4'	35:BA:1254:A:H1'	2.05	0.57
38:BD:142:VAL:HG23	38:BD:193:VAL:N	2.19	0.57
39:BE:199:ARG:HG3	39:BE:199:ARG:HH11	1.68	0.57
35:BA:2277:G:OP1	47:BQ:85:LYS:HB3	2.04	0.57
48:BR:41:ALA:O	48:BR:43:GLU:N	2.37	0.57
49:BS:49:VAL:HG21	49:BS:77:ALA:HB2	1.86	0.57
49:BS:92:TYR:HD1	49:BS:93:LYS:N	2.03	0.57
50:BT:92:GLY:C	50:BT:94:ALA:N	2.56	0.57
51:BU:95:LEU:HD12	52:BV:11:GLN:HG3	1.87	0.57
54:BX:39:ILE:C	54:BX:39:ILE:HD12	2.25	0.57
54:BX:60:ARG:HG3	54:BX:71:GLY:HA3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:11:ASP:N	55:BY:27:VAL:HG22	2.20	0.57
56:BZ:125:LEU:O	56:BZ:126:VAL:HG13	2.05	0.57
56:BZ:136:PHE:HD1	56:BZ:136:PHE:C	2.07	0.57
35:BA:875:G:O2'	56:BZ:151:HIS:NE2	2.33	0.57
56:BZ:167:PRO:O	56:BZ:168:GLU:HB2	2.05	0.57
1:CA:1018:C:H2'	1:CA:1019:C:C6	2.40	0.57
1:CA:1281:U:H5'	1:CA:1282:C:C5	2.40	0.57
1:CA:341:C:O2'	1:CA:342:C:H5'	2.05	0.57
1:CA:441:A:H3'	1:CA:442:C:C6	2.40	0.57
1:CA:962:C:H2'	1:CA:963:G:C8	2.38	0.57
1:CA:1073:U:O2	2:CB:104:ASN:ND2	2.38	0.57
2:CB:219:VAL:C	2:CB:222:ILE:HG22	2.25	0.57
2:CB:80:ILE:HG13	2:CB:81:VAL:N	2.20	0.57
7:CG:143:ARG:O	7:CG:147:ALA:HB2	2.05	0.57
9:CI:79:LEU:CD2	9:CI:102:LEU:HA	2.33	0.57
11:CK:109:VAL:HG22	18:CR:86:VAL:HA	1.86	0.57
11:CK:110:ASP:O	18:CR:84:LYS:HD2	2.05	0.57
14:CN:40:CYS:SG	14:CN:43:CYS:N	2.72	0.57
21:CU:17:THR:O	21:CU:22:ARG:HD3	2.05	0.57
25:CY:170:ALA:O	25:CY:171:LYS:C	2.43	0.57
31:D5:20:ARG:O	31:D5:21:SER:C	2.42	0.57
31:D5:31:VAL:HB	31:D5:32:PRO:CD	2.32	0.57
31:D5:44:THR:CG2	31:D5:45:VAL:H	2.11	0.57
34:D8:30:ARG:O	34:D8:30:ARG:HG3	2.04	0.57
35:DA:1151:G:H5''	51:DU:81:HIS:CE1	2.40	0.57
35:DA:1380:G:C2	35:DA:1381:G:C8	2.93	0.57
35:DA:1399:C:H2'	35:DA:1400:G:H8	1.68	0.57
35:DA:1528(A):A:H2'	35:DA:1529:G:H5''	1.87	0.57
35:DA:1876:A:H2'	35:DA:1877:A:H8	1.70	0.57
35:DA:2626:C:H2'	35:DA:2627:G:H8	1.70	0.57
35:DA:2802:G:O2'	35:DA:2803:C:C5'	2.52	0.57
35:DA:586:A:H2	35:DA:809:G:N3	2.03	0.57
38:DD:231:HIS:ND1	38:DD:232:PRO:CD	2.66	0.57
38:DD:268:ARG:HH12	38:DD:269:PHE:HE1	1.51	0.57
39:DE:38:THR:HG23	39:DE:39:PRO:HD2	1.86	0.57
40:DF:3:GLU:CG	40:DF:19:GLU:HB2	2.34	0.57
43:DI:10:GLU:O	43:DI:11:ASN:HB3	2.04	0.57
43:DI:72:LEU:O	43:DI:138:ILE:HG12	2.05	0.57
44:DN:10:GLU:HG3	44:DN:11:PRO:HD2	1.87	0.57
46:DP:45:LEU:HD23	46:DP:46:LYS:N	2.18	0.57
48:DR:37:THR:HG23	48:DR:40:LYS:HE2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:82:ILE:O	49:DS:83:LYS:HB2	2.04	0.57
1:AA:1202:G:H2'	1:AA:1203:C:C5'	2.33	0.57
1:AA:1305:G:H5'	21:AU:4:GLY:CA	2.35	0.57
1:AA:1396:A:C4'	1:AA:1398:A:H1'	2.35	0.57
1:AA:1459:C:O2'	1:AA:1460:A:H5'	2.05	0.57
2:AB:238:LEU:O	2:AB:239:VAL:C	2.44	0.57
3:AC:206:GLU:O	3:AC:208:ILE:N	2.37	0.57
3:AC:83:ARG:C	3:AC:87:LEU:HG	2.26	0.57
4:AD:62:GLN:HB3	4:AD:66:ARG:NH2	2.19	0.57
5:AE:10:MET:O	5:AE:10:MET:HG2	2.05	0.57
5:AE:35:GLY:HA2	5:AE:41:VAL:HG12	1.85	0.57
6:AF:58:GLY:O	6:AF:60:PHE:HD1	1.88	0.57
6:AF:52:ILE:O	6:AF:86:ARG:NH1	2.38	0.57
2:AB:178:ARG:NH2	8:AH:74:PRO:HG3	2.20	0.57
20:AT:14:LYS:HA	20:AT:17:ARG:HE	1.70	0.57
27:B1:38:SER:C	27:B1:39:LYS:HD3	2.26	0.57
32:B6:12:GLU:HB3	32:B6:23:THR:HG22	1.86	0.57
34:B8:39:LYS:HG2	34:B8:42:ARG:HH11	1.68	0.57
35:BA:1041:C:H5'	35:BA:1042:G:OP2	2.03	0.57
35:BA:1754:C:OP1	50:BT:96:ARG:NH1	2.36	0.57
35:BA:2182:G:H2'	35:BA:2183:C:C6	2.40	0.57
35:BA:375:C:H2'	35:BA:376:C:C6	2.40	0.57
35:BA:674:G:H1'	40:BF:74:ARG:HG3	1.87	0.57
36:BB:78:A:C2	36:BB:100:A:C4	2.93	0.57
38:BD:155:LEU:O	38:BD:156:ALA:C	2.42	0.57
38:BD:142:VAL:HG23	38:BD:192:THR:C	2.25	0.57
35:BA:1568:G:H4'	38:BD:59:LYS:CG	2.35	0.57
38:BD:25:THR:HG21	38:BD:82:ILE:N	2.19	0.57
40:BF:3:GLU:CG	40:BF:19:GLU:HB2	2.35	0.57
46:BP:138:LEU:CD2	46:BP:142:GLY:HA3	2.34	0.57
46:BP:7:ARG:HB3	46:BP:8:PRO:CD	2.32	0.57
48:BR:38:VAL:HB	48:BR:39:PRO:CD	2.30	0.57
50:BT:102:ILE:HG12	50:BT:103:ARG:N	2.20	0.57
50:BT:27:THR:O	50:BT:28:VAL:CB	2.52	0.57
50:BT:85:LYS:O	50:BT:85:LYS:HG2	2.04	0.57
54:BX:12:VAL:HG13	54:BX:17:ALA:CB	2.34	0.57
54:BX:28:PHE:N	54:BX:28:PHE:CD1	2.73	0.57
56:BZ:56:VAL:HA	56:BZ:70:LEU:HG	1.85	0.57
1:CA:989:C:N4	1:CA:1216:G:H1	2.01	0.57
1:CA:1283:G:O2'	1:CA:1284:C:H5'	2.04	0.57
1:CA:1294:G:H2'	1:CA:1295:G:C8	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:148:G:H2'	1:CA:149:A:C8	2.37	0.57
1:CA:673:G:H2'	1:CA:674:G:H8	1.68	0.57
2:CB:166:ASP:HB2	2:CB:205:ASP:OD2	2.04	0.57
2:CB:167:PRO:HG2	2:CB:192:SER:OG	2.04	0.57
2:CB:200:ILE:O	2:CB:201:ILE:HD13	2.04	0.57
2:CB:212:GLN:HG3	2:CB:235:SER:HB2	1.86	0.57
4:CD:100:ARG:CZ	4:CD:137:SER:HA	2.34	0.57
7:CG:75:VAL:HA	7:CG:88:PRO:HA	1.85	0.57
8:CH:96:GLY:O	8:CH:98:LYS:N	2.38	0.57
10:CJ:23:ILE:HG22	10:CJ:23:ILE:O	2.05	0.57
12:CL:6:THR:HG22	12:CL:9:GLN:CG	2.29	0.57
15:CO:33:THR:HG23	15:CO:63:ARG:NH1	2.20	0.57
17:CQ:67:LYS:CA	17:CQ:70:ARG:HH12	2.18	0.57
35:DA:1240:U:O2'	35:DA:1241:A:H5'	2.04	0.57
35:DA:1242:A:N1	46:DP:8:PRO:HG3	2.19	0.57
35:DA:1653:G:O2'	35:DA:1654:A:OP2	2.21	0.57
35:DA:1707:G:H2'	35:DA:1708:C:H6	1.69	0.57
35:DA:2650:U:O2'	35:DA:2651:C:H5'	2.05	0.57
38:DD:201:HIS:O	38:DD:203:ASN:N	2.37	0.57
39:DE:115:GLY:HA2	39:DE:157:ALA:HB1	1.86	0.57
39:DE:51:PHE:H	39:DE:74:PRO:CB	2.18	0.57
43:DI:92:VAL:HG22	43:DI:92:VAL:O	2.04	0.57
45:DO:89:ASN:C	45:DO:91:LEU:H	2.07	0.57
46:DP:122:PRO:CG	46:DP:141:ALA:HB3	2.34	0.57
47:DQ:69:PHE:CD1	47:DQ:70:PRO:HD2	2.40	0.57
48:DR:41:ALA:O	48:DR:44:LEU:N	2.38	0.57
48:DR:42:LYS:HG3	48:DR:45:ARG:NH2	2.20	0.57
39:DE:14:ILE:HB	50:DT:14:TYR:CE2	2.39	0.57
51:DU:44:ASN:O	51:DU:47:TYR:HB3	2.04	0.57
53:DW:74:ALA:C	53:DW:75:TYR:CD1	2.78	0.57
55:DY:77:PRO:O	55:DY:78:ALA:HB2	2.04	0.57
1:AA:962:C:H2'	1:AA:963:G:C8	2.39	0.56
3:AC:66:VAL:O	3:AC:66:VAL:HG12	2.04	0.56
3:AC:76:VAL:HG23	3:AC:77:ILE:N	2.20	0.56
8:AH:104:ARG:O	8:AH:107:LEU:HB3	2.05	0.56
8:AH:82:HIS:CD2	8:AH:138:TRP:NE1	2.69	0.56
16:AP:73:LEU:O	16:AP:77:ALA:HB2	2.04	0.56
18:AR:56:THR:OG1	18:AR:57:GLY:N	2.38	0.56
25:AY:130:ARG:HH11	25:AY:130:ARG:CG	2.17	0.56
25:AY:74:ASN:O	25:AY:77:LYS:HB2	2.05	0.56
35:BA:1642:G:H2'	35:BA:1643:G:C8	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:176:G:C2'	35:BA:177:G:H5'	2.34	0.56
35:BA:363(A):A:H2'	35:BA:363(A):A:N3	2.19	0.56
35:BA:768:G:H2'	35:BA:769:G:H8	1.70	0.56
36:BB:7:G:H4'	49:BS:29:PHE:HD2	1.70	0.56
38:BD:65:ILE:HD12	38:BD:65:ILE:O	2.05	0.56
41:BG:60:LEU:HD13	41:BG:61:ALA:N	2.20	0.56
43:BI:72:LEU:CD1	43:BI:138:ILE:HD11	2.23	0.56
43:BI:8:PRO:HD3	43:BI:15:VAL:CG1	2.35	0.56
43:BI:84:GLY:O	43:BI:85:GLU:HB2	2.05	0.56
44:BN:23:LEU:HB3	44:BN:60:ILE:HG21	1.87	0.56
44:BN:40:PRO:HG3	51:BU:68:ALA:HB2	1.87	0.56
44:BN:78:TYR:N	44:BN:79:PRO:CD	2.68	0.56
45:BO:14:THR:HG22	45:BO:52:VAL:HG21	1.85	0.56
46:BP:80:TYR:CD1	46:BP:111:ARG:HB3	2.41	0.56
47:BQ:109:VAL:HG13	47:BQ:113:GLN:OE1	2.05	0.56
49:BS:85:VAL:HG23	49:BS:106:ARG:HB2	1.87	0.56
50:BT:109:GLU:O	50:BT:112:ARG:HG3	2.04	0.56
54:BX:72:LYS:HE3	54:BX:74:PRO:CB	2.27	0.56
1:CA:1261:A:H62	1:CA:1274:G:H21	1.53	0.56
1:CA:1478:C:H2'	1:CA:1479:C:H6	1.68	0.56
1:CA:622:A:C8	1:CA:623:C:C5	2.92	0.56
1:CA:735:C:H2'	1:CA:736:C:C6	2.36	0.56
1:CA:802:A:H2'	1:CA:803:G:C5'	2.34	0.56
1:CA:949:A:H2'	1:CA:950:U:O4'	2.04	0.56
2:CB:63:MET:HB3	2:CB:225:ALA:HB1	1.86	0.56
3:CC:113:ALA:HB3	3:CC:114:PRO:HD3	1.87	0.56
3:CC:47:LEU:HD21	3:CC:52:LEU:HD13	1.87	0.56
6:CF:29:ALA:HB1	6:CF:79:LEU:CD2	2.35	0.56
7:CG:26:PHE:HB2	7:CG:62:PHE:CZ	2.39	0.56
9:CI:17:VAL:CG2	9:CI:81:ILE:HD13	2.34	0.56
11:CK:49:GLY:O	11:CK:50:TYR:HD2	1.88	0.56
13:CM:68:GLY:O	13:CM:69:GLU:HB2	2.04	0.56
18:CR:22:VAL:HA	18:CR:25:THR:HG1	1.70	0.56
18:CR:44:LEU:HA	18:CR:49:LYS:O	2.05	0.56
19:CS:14:HIS:CD2	19:CS:15:LEU:HD22	2.40	0.56
27:D1:34:THR:HG21	35:DA:388:G:P	2.45	0.56
35:DA:1002:G:H2'	35:DA:1003:G:O4'	2.05	0.56
35:DA:1427:A:H4'	35:DA:1428:C:O5'	2.03	0.56
35:DA:1935:G:H1'	35:DA:1964:G:N2	2.20	0.56
35:DA:2086:U:OP1	38:DD:262:ARG:HG2	2.05	0.56
35:DA:534:U:O3'	51:DU:46:ALA:HB2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DC:55:ASP:CG	37:DC:56:GLN:H	2.09	0.56
39:DE:103:ASP:OD1	39:DE:168:MET:HG2	2.05	0.56
35:DA:2680:C:H5'	39:DE:189:PRO:HA	1.87	0.56
40:DF:192:LEU:HD21	40:DF:194:MET:HE2	1.87	0.56
40:DF:71:GLY:O	40:DF:72:ARG:C	2.42	0.56
42:DH:150:ALA:O	42:DH:152:ARG:N	2.37	0.56
42:DH:89:ILE:H	42:DH:89:ILE:HD13	1.69	0.56
50:DT:35:LYS:HE2	50:DT:41:ARG:HG3	1.87	0.56
50:DT:77:PRO:O	50:DT:78:LEU:HB3	2.06	0.56
51:DU:92:ARG:HH22	52:DV:10:LYS:HA	1.69	0.56
52:DV:19:LYS:HG3	52:DV:20:LEU:H	1.68	0.56
52:DV:40:LEU:O	52:DV:41:GLY:C	2.42	0.56
54:DX:77:LYS:CD	54:DX:78:LYS:HG3	2.35	0.56
56:DZ:24:LEU:HD12	56:DZ:25:PRO:N	2.20	0.56
1:AA:1097:C:O2'	1:AA:1098:C:H5'	2.05	0.56
1:AA:1321:C:H5''	1:AA:1322:C:H2'	1.86	0.56
1:AA:1486:G:H2'	1:AA:1487:G:O4'	2.05	0.56
1:AA:35:G:H2'	1:AA:36:C:C6	2.40	0.56
1:AA:690:G:H2'	1:AA:691:G:O4'	2.05	0.56
1:AA:826:C:H2'	1:AA:827:U:C6	2.40	0.56
2:AB:50:GLU:OE1	2:AB:200:ILE:HB	2.05	0.56
4:AD:30:LYS:HB3	4:AD:35:ARG:HH11	1.70	0.56
5:AE:12:LEU:CD1	5:AE:31:LEU:HB3	2.35	0.56
8:AH:112:LEU:C	8:AH:112:LEU:HD12	2.25	0.56
8:AH:83:ILE:CB	8:AH:137:VAL:HG13	2.34	0.56
18:AR:86:VAL:HG12	18:AR:87:ARG:NH1	2.20	0.56
20:AT:8:ARG:HD2	20:AT:8:ARG:N	2.18	0.56
23:AW:27:G:H2'	23:AW:28:U:H6	1.69	0.56
25:AY:25:LEU:O	25:AY:28:LEU:HB2	2.05	0.56
35:BA:1298:C:H3'	35:BA:1299:G:H8	1.69	0.56
35:BA:1405:U:H2'	35:BA:1406:U:H6	1.68	0.56
35:BA:1440:G:H2'	35:BA:1441:G:H8	1.70	0.56
35:BA:1528(A):A:C2'	35:BA:1529:G:H5''	2.35	0.56
35:BA:1826:G:H4'	38:BD:242:ARG:NH2	2.19	0.56
35:BA:1930:G:H22	35:BA:1968:G:C2'	2.16	0.56
35:BA:2055:C:H4'	35:BA:2056:G:H5''	1.87	0.56
35:BA:2056:G:H2'	35:BA:2056:G:N3	2.20	0.56
35:BA:2187:G:O2'	35:BA:2188:C:H5'	2.06	0.56
35:BA:2801(A):A:H5'	35:BA:2802:G:C8	2.40	0.56
35:BA:669:G:N3	35:BA:669:G:H2'	2.19	0.56
37:BC:214:VAL:C	37:BC:216:THR:H	2.08	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:103:ASP:OD2	39:BE:202:LYS:HE2	2.05	0.56
41:BG:76:SER:HA	41:BG:83:ARG:HA	1.86	0.56
43:BI:130:TYR:CB	43:BI:136:VAL:HG13	2.35	0.56
43:BI:82:ARG:HG3	43:BI:82:ARG:HH11	1.71	0.56
43:BI:83:ALA:HA	43:BI:89:TYR:CD1	2.39	0.56
47:BQ:141:GLN:HA	56:BZ:71:VAL:O	2.05	0.56
52:BV:61:VAL:HG23	52:BV:100:ARG:N	2.20	0.56
55:BY:7:VAL:CG2	55:BY:8:LYS:HD2	2.35	0.56
1:CA:1010:G:N2	1:CA:1020:U:H1'	2.20	0.56
1:CA:1223:C:P	19:CS:78:ARG:HH22	2.28	0.56
1:CA:501:C:H2'	1:CA:502:G:H8	1.69	0.56
1:CA:637:G:H2'	1:CA:638:G:H8	1.70	0.56
1:CA:893:C:H2'	1:CA:894:G:C8	2.40	0.56
4:CD:116:GLN:NE2	4:CD:157:LEU:HD21	2.18	0.56
10:CJ:48:THR:HG23	10:CJ:61:GLU:C	2.26	0.56
10:CJ:6:ILE:O	10:CJ:71:LEU:HD12	2.04	0.56
11:CK:28:THR:HG22	11:CK:29:ILE:N	2.19	0.56
13:CM:84:ILE:HG22	13:CM:84:ILE:O	2.05	0.56
15:CO:53:HIS:CE1	15:CO:57:LEU:HD21	2.40	0.56
31:D5:26:THR:O	31:D5:26:THR:HG23	2.05	0.56
34:D8:35:GLN:HE21	34:D8:36:LYS:NZ	2.02	0.56
34:D8:56:GLU:C	34:D8:58:ILE:N	2.57	0.56
34:D8:59:LYS:C	34:D8:61:LEU:N	2.45	0.56
35:DA:1131:G:N3	35:DA:1132:A:C8	2.74	0.56
35:DA:1274:A:N3	35:DA:1297:C:H1'	2.20	0.56
35:DA:1528:A:H2'	35:DA:1528:A:N3	2.19	0.56
35:DA:1710:C:H2'	35:DA:1711:C:C6	2.40	0.56
35:DA:1857:G:O2'	35:DA:1885:A:N6	2.39	0.56
35:DA:2339:G:H2'	35:DA:2340:G:H8	1.70	0.56
35:DA:236:C:O2'	35:DA:237:C:H5'	2.06	0.56
35:DA:2666:C:H2'	35:DA:2667:C:O4'	2.05	0.56
35:DA:532:A:H2'	35:DA:532:A:N3	2.19	0.56
35:DA:916:G:C2'	35:DA:917:A:H5''	2.36	0.56
35:DA:950:G:O2'	35:DA:951:C:H5'	2.04	0.56
36:DB:41:U:C4	41:DG:70:VAL:O	2.58	0.56
44:DN:128:HIS:O	44:DN:130:HIS:N	2.38	0.56
44:DN:78:TYR:N	44:DN:79:PRO:CD	2.69	0.56
51:DU:91:ASP:O	51:DU:92:ARG:O	2.23	0.56
51:DU:99:ALA:HB2	51:DU:106:PHE:CD1	2.40	0.56
53:DW:75:TYR:HE1	53:DW:104:THR:HB	1.70	0.56
1:AA:103:C:H2'	1:AA:104:G:H8	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1269:A:H2	1:AA:1312:G:N3	2.03	0.56
1:AA:453:A:C5	1:AA:454:C:C4	2.94	0.56
1:AA:884:U:H4'	1:AA:885:G:C5'	2.34	0.56
3:AC:76:VAL:HG23	3:AC:77:ILE:HG13	1.86	0.56
4:AD:96:LEU:H	4:AD:96:LEU:CD2	1.97	0.56
9:AI:7:THR:HB	9:AI:83:ARG:HH11	1.70	0.56
18:AR:70:ILE:O	18:AR:74:ARG:HB2	2.05	0.56
1:AA:1223:C:P	19:AS:78:ARG:HH22	2.28	0.56
20:AT:32:ALA:O	20:AT:33:ILE:C	2.42	0.56
25:AY:103:ILE:CG1	25:AY:103:ILE:O	2.53	0.56
27:B1:33:LYS:CG	27:B1:34:THR:H	2.16	0.56
28:B2:46:GLN:NE2	28:B2:47:ASN:N	2.53	0.56
35:BA:1240:U:O2'	35:BA:1241:A:H5'	2.04	0.56
35:BA:137:C:C2'	35:BA:139:G:H5'	2.35	0.56
35:BA:1655:A:H4'	39:BE:115:GLY:H	1.69	0.56
35:BA:1819:A:O2'	35:BA:1820:U:OP2	2.19	0.56
35:BA:2266:A:H4'	35:BA:2267:A:C2	2.41	0.56
35:BA:271(J):C:C2'	35:BA:271(K):U:H5''	2.36	0.56
35:BA:271(J):C:C3'	35:BA:271(K):U:H5''	2.35	0.56
35:BA:549:G:H2'	35:BA:551:G:O4'	2.06	0.56
35:BA:797:C:OP2	40:BF:62:ARG:HG3	2.06	0.56
35:BA:816:C:O2'	35:BA:817:C:H5'	2.05	0.56
35:BA:878:A:H3'	35:BA:879:G:C8	2.39	0.56
35:BA:972:G:H2'	35:BA:973:A:C8	2.40	0.56
38:BD:108:PRO:HG2	38:BD:111:LEU:HB2	1.88	0.56
38:BD:125:ILE:H	38:BD:125:ILE:CD1	2.19	0.56
35:BA:2203:U:C4'	38:BD:151:LYS:HE3	2.35	0.56
40:BF:170:LEU:HD21	40:BF:172:TRP:CE2	2.40	0.56
40:BF:81:PRO:HG2	40:BF:82:ILE:H	1.70	0.56
41:BG:165:THR:HB	41:BG:167:GLU:OE1	2.05	0.56
45:BO:22:ILE:H	45:BO:41:ALA:HA	1.70	0.56
46:BP:123:LEU:O	46:BP:125:VAL:HG12	2.05	0.56
35:BA:1242:A:N1	46:BP:8:PRO:HG3	2.20	0.56
49:BS:80:LEU:HD12	49:BS:80:LEU:N	2.20	0.56
52:BV:34:GLU:HG2	52:BV:35:LEU:N	2.20	0.56
53:BW:8:ARG:HH11	53:BW:8:ARG:HG3	1.71	0.56
35:BA:1341:U:O3'	54:BX:55:ASN:HB3	2.05	0.56
55:BY:6:HIS:ND1	55:BY:6:HIS:N	2.50	0.56
1:CA:1074:G:H4'	2:CB:103:THR:HG22	1.87	0.56
1:CA:1154:G:H2'	1:CA:1155:G:H8	1.70	0.56
1:CA:1296:C:H5'	1:CA:1297:C:OP2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1360:A:H2'	1:CA:1361:G:O4'	2.05	0.56
1:CA:20:U:O2'	1:CA:21:G:H5'	2.05	0.56
1:CA:552:U:O3'	12:CL:87:GLY:HA3	2.06	0.56
1:CA:685:G:N2	1:CA:706:A:H61	2.03	0.56
2:CB:159:PRO:C	2:CB:161:ALA:N	2.56	0.56
2:CB:36:ARG:NE	2:CB:37:ASN:H	2.03	0.56
11:CK:21:ILE:CG1	11:CK:84:VAL:HG12	2.35	0.56
15:CO:56:LEU:HA	15:CO:59:MET:HE2	1.87	0.56
16:CP:6:LEU:HB3	16:CP:17:TYR:HB3	1.86	0.56
18:CR:36:ASN:HD22	18:CR:39:VAL:HB	1.68	0.56
19:CS:12:ASP:O	19:CS:16:LEU:HD13	2.04	0.56
27:D1:18:ILE:HA	27:D1:44:PRO:HD2	1.87	0.56
35:DA:1657:C:O2'	35:DA:1658:C:H5'	2.05	0.56
35:DA:1658:C:OP1	39:DE:132:HIS:ND1	2.38	0.56
35:DA:1788:C:H2'	35:DA:1789:A:C8	2.34	0.56
35:DA:1841:U:H2'	35:DA:1842:G:H8	1.71	0.56
35:DA:2019:A:H4'	51:DU:34:LYS:HD2	1.85	0.56
35:DA:2081:C:H2'	35:DA:2082:A:H8	1.69	0.56
35:DA:2579:C:H4'	39:DE:134:ILE:HG13	1.87	0.56
35:DA:2773:C:H2'	35:DA:2774:C:C6	2.40	0.56
35:DA:985:C:H2'	35:DA:985:C:O2	2.06	0.56
37:DC:214:VAL:C	37:DC:216:THR:H	2.09	0.56
38:DD:131:LEU:N	38:DD:131:LEU:HD12	2.20	0.56
42:DH:138:LYS:H	42:DH:141:VAL:HG23	1.68	0.56
44:DN:53:VAL:HG13	44:DN:121:LYS:O	2.05	0.56
45:DO:63:VAL:HG22	45:DO:83:ALA:O	2.06	0.56
45:DO:87:ILE:HG23	45:DO:88:ASN:N	2.20	0.56
47:DQ:8:LYS:CG	47:DQ:9:TYR:N	2.67	0.56
49:DS:80:LEU:HD12	49:DS:80:LEU:N	2.20	0.56
50:DT:96:ARG:CG	50:DT:96:ARG:HH11	2.17	0.56
35:DA:445:C:O3'	51:DU:3:ARG:HG2	2.04	0.56
51:DU:92:ARG:NH1	52:DV:11:GLN:O	2.38	0.56
51:DU:90:VAL:HG22	52:DV:39:LEU:CD1	2.35	0.56
53:DW:17:VAL:O	53:DW:20:VAL:HG23	2.06	0.56
54:DX:21:PHE:CE1	54:DX:26:TYR:HB3	2.39	0.56
56:DZ:50:GLN:OE1	56:DZ:50:GLN:N	2.37	0.56
1:AA:10:A:H2'	1:AA:11:G:H8	1.71	0.56
1:AA:1283:G:O2'	1:AA:1284:C:H5'	2.04	0.56
1:AA:1296:C:H3'	1:AA:1297:C:H6	1.71	0.56
1:AA:1392:G:N2	1:AA:1502:A:C8	2.72	0.56
1:AA:391:G:N1	1:AA:392:G:C5	2.73	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:441:A:H3'	1:AA:442:C:H6	1.69	0.56
1:AA:126:G:H5'	1:AA:633:G:N2	2.19	0.56
1:AA:64:G:OP1	1:AA:64:G:H3'	2.05	0.56
1:AA:732:C:H2'	1:AA:733:A:H5''	1.87	0.56
2:AB:237:ALA:O	2:AB:238:LEU:HB3	2.04	0.56
4:AD:10:ARG:O	4:AD:13:ARG:HB3	2.04	0.56
7:AG:71:PRO:CG	7:AG:103:TRP:HZ3	2.18	0.56
12:AL:89:ARG:HH11	12:AL:89:ARG:HB2	1.69	0.56
18:AR:36:ASN:HD22	18:AR:39:VAL:HB	1.70	0.56
19:AS:6:LYS:HD2	19:AS:7:LYS:HD2	1.86	0.56
27:B1:11:ARG:NH1	27:B1:59:THR:O	2.37	0.56
27:B1:76:ARG:CA	27:B1:78:LYS:HZ3	2.18	0.56
29:B3:41:PRO:HD3	29:B3:44:ARG:CZ	2.35	0.56
31:B5:43:HIS:CD2	31:B5:43:HIS:N	2.72	0.56
35:BA:1528:A:H2'	35:BA:1528:A:N3	2.19	0.56
35:BA:2114:A:C2'	35:BA:2115:G:H5'	2.34	0.56
35:BA:2340:G:O2'	35:BA:2341:G:H5'	2.05	0.56
35:BA:826:U:OP1	35:BA:2428:G:OP1	2.23	0.56
35:BA:2864:G:O2'	35:BA:2865:U:H5'	2.05	0.56
33:B7:16:HIS:CE1	35:BA:465:G:H4'	2.40	0.56
35:BA:619:G:P	35:BA:620:G:H22	2.28	0.56
35:BA:779:U:H2'	35:BA:780:G:C8	2.41	0.56
38:BD:120:GLY:O	38:BD:131:LEU:HB3	2.05	0.56
38:BD:94:LEU:HA	38:BD:104:TYR:HA	1.87	0.56
43:BI:10:GLU:O	43:BI:12:LEU:HD23	2.05	0.56
43:BI:4:ILE:C	43:BI:5:LEU:HD23	2.26	0.56
44:BN:4:TYR:CD1	44:BN:4:TYR:N	2.71	0.56
45:BO:32:TYR:N	45:BO:32:TYR:HD1	2.02	0.56
45:BO:43:VAL:O	45:BO:45:GLU:N	2.37	0.56
35:BA:943:U:OP1	46:BP:38:GLN:HB3	2.04	0.56
51:BU:26:GLY:C	51:BU:28:ARG:N	2.57	0.56
51:BU:90:VAL:CG1	52:BV:39:LEU:HB3	2.33	0.56
53:BW:11:ARG:HH22	53:BW:98:LYS:HB3	1.70	0.56
54:BX:29:TRP:CZ3	54:BX:76:ARG:HG2	2.38	0.56
1:CA:1411:C:O2'	1:CA:1412:C:H5'	2.05	0.56
1:CA:1504:G:H4'	1:CA:1505:G:C4	2.40	0.56
1:CA:361:G:O2'	1:CA:362:G:H5'	2.06	0.56
1:CA:950:U:H2'	1:CA:951:G:H8	1.71	0.56
4:CD:119:GLN:HB3	4:CD:120:LEU:HD12	1.87	0.56
5:CE:129:ILE:O	5:CE:132:ALA:N	2.37	0.56
7:CG:21:VAL:HG23	7:CG:22:LEU:H	1.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:84:ASN:N	7:CG:84:ASN:ND2	2.54	0.56
15:CO:15:PHE:O	15:CO:16:ALA:O	2.24	0.56
15:CO:39:LEU:HD12	15:CO:56:LEU:CD1	2.36	0.56
17:CQ:48:GLU:O	17:CQ:50:LYS:N	2.38	0.56
20:CT:26:ASN:ND2	20:CT:26:ASN:N	2.54	0.56
20:CT:32:ALA:O	20:CT:33:ILE:C	2.43	0.56
20:CT:32:ALA:O	20:CT:36:LEU:HD23	2.05	0.56
20:CT:41:ILE:O	20:CT:43:LEU:N	2.39	0.56
25:CY:122:ALA:O	25:CY:123:GLU:C	2.44	0.56
25:CY:161:ILE:O	25:CY:164:ILE:HB	2.05	0.56
25:CY:23:HIS:O	25:CY:25:LEU:N	2.38	0.56
25:CY:43:VAL:HG23	25:CY:50:VAL:O	2.04	0.56
35:DA:1430:C:H2'	35:DA:1431:U:C6	2.40	0.56
35:DA:1821:A:H2'	35:DA:1822:G:C8	2.39	0.56
35:DA:2266:A:H4'	35:DA:2267:A:N3	2.20	0.56
29:D3:14:GLY:O	35:DA:969:U:H4'	2.05	0.56
35:DA:2729:G:C1'	39:DE:187:ALA:HB2	2.23	0.56
41:DG:35:GLU:HG2	41:DG:35:GLU:O	2.06	0.56
42:DH:149:ARG:CG	42:DH:162:ILE:HD11	2.35	0.56
47:DQ:29:PHE:CD1	47:DQ:29:PHE:N	2.72	0.56
51:DU:33:ARG:HA	51:DU:36:ARG:HB2	1.86	0.56
51:DU:47:TYR:HA	51:DU:50:ARG:NH2	2.20	0.56
54:DX:53:LYS:HZ2	54:DX:55:ASN:HD21	1.52	0.56
54:DX:88:LYS:C	54:DX:90:GLU:H	2.07	0.56
55:DY:74:PRO:HG2	55:DY:80:GLY:O	2.06	0.56
55:DY:83:THR:O	55:DY:84:ARG:HG3	2.05	0.56
1:AA:1096:C:O2'	1:AA:1097:C:H5'	2.05	0.56
1:AA:955:U:C1'	1:AA:1227:A:H61	2.15	0.56
1:AA:1405:G:H2'	1:AA:1406:U:H6	1.70	0.56
1:AA:158:G:O2'	1:AA:159:G:H5'	2.05	0.56
1:AA:191:G:C4	20:AT:105:SER:HB3	2.41	0.56
1:AA:33:A:H2'	1:AA:34:C:C6	2.40	0.56
1:AA:452:A:C2	1:AA:453:A:C4	2.93	0.56
1:AA:766:A:H2'	1:AA:767:A:O4'	2.05	0.56
4:AD:108:LEU:O	4:AD:110:PHE:CD1	2.58	0.56
9:AI:102:LEU:O	9:AI:103:THR:OG1	2.21	0.56
10:AJ:4:ILE:CB	10:AJ:74:ILE:HD11	2.28	0.56
11:AK:86:GLY:H	11:AK:112:THR:CG2	2.19	0.56
12:AL:119:LYS:C	12:AL:120:TYR:HD1	2.07	0.56
18:AR:75:ILE:C	18:AR:76:LEU:HD22	2.26	0.56
26:B0:1:MET:O	26:B0:2:ALA:HB3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:B0:21:LEU:CD1	26:B0:41:ARG:HD3	2.35	0.56
27:B1:11:ARG:HG3	27:B1:61:ARG:C	2.25	0.56
27:B1:51:VAL:CG2	27:B1:62:VAL:HG11	2.34	0.56
31:B5:57:VAL:HG23	31:B5:58:LEU:N	2.18	0.56
34:B8:35:GLN:HE21	34:B8:36:LYS:NZ	2.03	0.56
34:B8:56:GLU:O	34:B8:59:LYS:NZ	2.36	0.56
35:BA:443:A:H1'	35:BA:1201:C:O4'	2.06	0.56
35:BA:1657:C:O2'	35:BA:1658:C:H5'	2.05	0.56
35:BA:1857:G:O2'	35:BA:1885:A:N6	2.39	0.56
35:BA:2010:G:H5''	53:BW:42:ARG:HB2	1.86	0.56
35:BA:2246:G:H2'	35:BA:2247:A:C8	2.40	0.56
35:BA:2290:G:H2'	35:BA:2291:U:C6	2.41	0.56
35:BA:2753:A:H2	35:BA:2754:U:C2	2.23	0.56
35:BA:745:G:H2'	35:BA:746:A:H5'	1.87	0.56
36:BB:42:C:O2	41:BG:93:THR:N	2.38	0.56
38:BD:260:ARG:NH1	38:BD:260:ARG:HG2	2.18	0.56
38:BD:96:HIS:CE1	38:BD:102:LYS:HD2	2.41	0.56
39:BE:63:LEU:O	39:BE:64:LYS:C	2.44	0.56
40:BF:41:LEU:HD23	40:BF:44:ARG:HD3	1.87	0.56
41:BG:144:ILE:HD12	41:BG:145:THR:H	1.71	0.56
41:BG:85:GLY:O	41:BG:86:MET:HB2	2.06	0.56
42:BH:102:ALA:CB	42:BH:117:PRO:HD3	2.20	0.56
45:BO:35:VAL:CG2	45:BO:69:ILE:HG12	2.35	0.56
40:BF:34:TRP:CZ3	46:BP:12:ALA:HA	2.40	0.56
46:BP:62:LEU:HD22	46:BP:62:LEU:C	2.25	0.56
47:BQ:42:ILE:HD13	47:BQ:97:VAL:CG2	2.35	0.56
49:BS:67:ARG:HE	49:BS:100:ALA:HB3	1.69	0.56
50:BT:106:SER:O	50:BT:107:ASP:HB3	2.06	0.56
50:BT:77:PRO:O	50:BT:78:LEU:HB3	2.06	0.56
52:BV:5:VAL:CG2	52:BV:36:PRO:HB2	2.33	0.56
54:BX:77:LYS:CD	54:BX:78:LYS:HG3	2.34	0.56
56:BZ:127:LYS:O	56:BZ:128:VAL:HB	2.04	0.56
1:CA:9:G:H2'	1:CA:10:A:H8	1.71	0.56
2:CB:137:ARG:NH1	2:CB:137:ARG:HG2	2.20	0.56
2:CB:164:VAL:O	2:CB:186:ALA:HB1	2.05	0.56
2:CB:34:ALA:O	2:CB:41:ILE:HB	2.06	0.56
3:CC:155:GLY:O	3:CC:156:ARG:HB2	2.05	0.56
6:CF:41:GLU:H	6:CF:62:TRP:HE3	1.52	0.56
7:CG:42:ILE:HG23	7:CG:117:ALA:HB2	1.87	0.56
8:CH:12:ARG:HH11	8:CH:26:VAL:HA	1.70	0.56
13:CM:13:LYS:HZ2	13:CM:21:TYR:HE1	1.52	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CT:50:GLU:HG3	20:CT:51:GLU:H	1.70	0.56
23:CW:75:C:O2'	27:D1:32:LYS:HD2	2.06	0.56
35:DA:1159:U:C2'	35:DA:1160:G:H5'	2.34	0.56
35:DA:1197:G:H5'	35:DA:1227:G:O2'	2.05	0.56
35:DA:1411:C:HO2'	35:DA:1412:A:H8	1.50	0.56
35:DA:1665:A:O2'	35:DA:1666:G:H5'	2.04	0.56
35:DA:2392:A:H1'	46:DP:60:MET:HE3	1.87	0.56
35:DA:2408:U:H2'	35:DA:2409:G:H8	1.68	0.56
35:DA:2574:G:H2'	35:DA:2575:C:H6	1.70	0.56
35:DA:28:A:H1'	35:DA:513:A:C2	2.40	0.56
35:DA:523:C:C2'	35:DA:524:U:H5'	2.36	0.56
39:DE:3:GLY:HA3	39:DE:81:ILE:HG21	1.88	0.56
40:DF:57:VAL:HG12	40:DF:58:ALA:N	2.19	0.56
43:DI:15:VAL:O	43:DI:16:GLY:C	2.43	0.56
46:DP:48:PRO:O	46:DP:51:PHE:N	2.38	0.56
54:DX:7:VAL:O	54:DX:31:HIS:N	2.35	0.56
55:DY:37:VAL:HG13	55:DY:69:ALA:CB	2.36	0.56
47:DQ:141:GLN:OE1	56:DZ:70:LEU:HB2	2.04	0.56
56:DZ:96:VAL:HG22	56:DZ:97:GLU:N	2.20	0.56
1:AA:1003:G:H2'	1:AA:1004:A:C4'	2.35	0.56
1:AA:1256:A:H61	1:AA:1278:U:H1'	1.69	0.56
1:AA:1281:U:H5'	1:AA:1282:C:C5	2.40	0.56
1:AA:259:G:H2'	1:AA:260:G:C8	2.41	0.56
1:AA:775:G:O2'	1:AA:776:G:H5'	2.06	0.56
1:AA:949:A:H61	1:AA:1232:U:H3	1.53	0.56
1:AA:1073:U:O2	2:AB:104:ASN:ND2	2.39	0.56
3:AC:155:GLY:O	3:AC:156:ARG:HB2	2.05	0.56
4:AD:133:VAL:CG1	4:AD:135:LEU:H	2.19	0.56
8:AH:6:ILE:O	8:AH:10:LEU:HG	2.06	0.56
9:AI:73:GLN:O	9:AI:77:ILE:HG13	2.05	0.56
29:B3:26:LEU:N	29:B3:26:LEU:HD23	2.20	0.56
34:B8:53:PRO:C	34:B8:55:ALA:H	2.07	0.56
35:BA:1414:G:H2'	35:BA:1415:U:C5	2.41	0.56
35:BA:1475:G:H2'	35:BA:1475:G:N3	2.20	0.56
35:BA:147:U:H2'	35:BA:148:C:H6	1.71	0.56
35:BA:1703:G:H2'	35:BA:1704:G:H8	1.70	0.56
35:BA:1839:G:H5'	35:BA:1839:G:H8	1.70	0.56
35:BA:1899:G:H22	35:BA:1902:C:N4	2.02	0.56
35:BA:2007:C:H2'	35:BA:2008:C:C6	2.37	0.56
35:BA:2650:U:O2'	35:BA:2651:C:H5'	2.06	0.56
35:BA:2711:A:OP1	35:BA:2712(A):A:P	2.63	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2821:A:H2'	35:BA:2822:G:H8	1.67	0.56
36:BB:21:G:N3	36:BB:21:G:H2'	2.20	0.56
37:BC:41:VAL:HB	37:BC:178:ALA:HB1	1.88	0.56
39:BE:59:VAL:HG22	39:BE:63:LEU:HA	1.87	0.56
44:BN:55:VAL:HG12	44:BN:56:ASN:N	2.21	0.56
47:BQ:124:LYS:HA	47:BQ:124:LYS:HE2	1.87	0.56
48:BR:82:GLU:OE1	48:BR:83:ILE:HD13	2.06	0.56
35:BA:2334:G:C5'	49:BS:13:ARG:HG2	2.30	0.56
49:BS:57:LYS:O	49:BS:58:LEU:HB2	2.06	0.56
49:BS:82:ILE:O	49:BS:83:LYS:HB2	2.03	0.56
49:BS:83:LYS:HE3	49:BS:84:GLN:HG3	1.86	0.56
55:BY:44:ILE:CG2	55:BY:45:VAL:N	2.69	0.56
56:BZ:129:SER:HB3	56:BZ:132:ASN:HD22	1.69	0.56
1:CA:1003:G:H2'	1:CA:1004:A:C4'	2.35	0.56
1:CA:1288:A:O2'	1:CA:1289:A:H5'	2.05	0.56
1:CA:1346:A:H61	1:CA:1374:A:H3'	1.70	0.56
1:CA:1442(B):A:OP1	1:CA:1442(B):A:H4'	2.05	0.56
1:CA:1452:C:H1'	1:CA:1456:G:C2	2.38	0.56
1:CA:358:U:H2'	1:CA:359:U:H6	1.71	0.56
1:CA:441:A:H3'	1:CA:442:C:H6	1.70	0.56
1:CA:811:C:H4'	1:CA:900:A:N6	2.20	0.56
1:CA:811:C:O2'	1:CA:901:A:N1	2.38	0.56
2:CB:99:GLY:O	2:CB:101:MET:N	2.39	0.56
2:CB:213:LEU:C	2:CB:213:LEU:HD23	2.25	0.56
2:CB:88:ALA:HB2	2:CB:223:ILE:HD11	1.86	0.56
6:CF:58:GLY:O	6:CF:60:PHE:HD1	1.89	0.56
8:CH:89:PRO:HA	8:CH:92:ARG:NH1	2.20	0.56
20:CT:19:SER:O	20:CT:23:ARG:N	2.36	0.56
26:D0:38:VAL:HG23	26:D0:59:LEU:HB2	1.85	0.56
27:D1:10:LYS:O	27:D1:11:ARG:O	2.23	0.56
29:D3:43:ILE:O	29:D3:47:VAL:HG23	2.05	0.56
35:DA:2024:G:H2'	35:DA:2025:C:H6	1.70	0.56
35:DA:2250:G:C8	35:DA:2496:C:H5''	2.40	0.56
35:DA:2636:U:H4'	39:DE:80:GLU:OE1	2.04	0.56
35:DA:565:C:H4'	35:DA:1253:A:N6	2.21	0.56
35:DA:614(C):A:N3	40:DF:180:GLY:HA2	2.20	0.56
35:DA:632:A:N3	35:DA:2403:C:H1'	2.21	0.56
35:DA:827:U:H2'	35:DA:2068:U:C2	2.40	0.56
35:DA:859:G:H5'	35:DA:2268:A:O2'	2.05	0.56
35:DA:85:G:O5'	55:DY:30:VAL:HB	2.06	0.56
35:DA:956:G:H5'	35:DA:957:A:OP2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:108:PRO:HB3	38:DD:143:HIS:CE1	2.41	0.56
38:DD:70:TRP:CH2	38:DD:150:LYS:HA	2.40	0.56
38:DD:231:HIS:CE1	38:DD:232:PRO:HD2	2.41	0.56
39:DE:30:PRO:C	39:DE:32:PRO:HD3	2.25	0.56
39:DE:55:ASN:ND2	39:DE:75:VAL:HG13	2.21	0.56
41:DG:102:PHE:CE2	41:DG:141:PHE:HE1	2.24	0.56
42:DH:35:VAL:HG12	42:DH:35:VAL:O	2.06	0.56
43:DI:68:LEU:HG	43:DI:72:LEU:HD23	1.86	0.56
35:DA:906:G:H5'	47:DQ:26:TYR:OH	2.05	0.56
47:DQ:42:ILE:HD13	47:DQ:97:VAL:CG2	2.35	0.56
47:DQ:20:ALA:CA	47:DQ:98:LYS:HB3	2.33	0.56
48:DR:2:ARG:NE	48:DR:5:LYS:HE3	2.21	0.56
50:DT:109:GLU:O	50:DT:112:ARG:HG3	2.06	0.56
56:DZ:48:PHE:HE2	56:DZ:71:VAL:CG2	2.12	0.56
56:DZ:56:VAL:HG12	56:DZ:57:ILE:N	2.21	0.56
1:AA:431:A:H2'	1:AA:432:A:H8	1.71	0.56
1:AA:519:C:O2'	1:AA:520:A:H5'	2.05	0.56
3:AC:130:VAL:HA	3:AC:133:ALA:HB3	1.86	0.56
7:AG:50:ILE:O	7:AG:54:THR:O	2.23	0.56
8:AH:20:TYR:HE2	8:AH:75:ARG:HD2	1.70	0.56
9:AI:14:VAL:O	9:AI:65:VAL:HG23	2.05	0.56
13:AM:84:ILE:O	13:AM:84:ILE:HG22	2.05	0.56
20:AT:58:LYS:HE3	20:AT:62:LEU:CD1	2.36	0.56
23:AW:36:A:O2'	23:AW:37:U:H5'	2.05	0.56
27:B1:86:SER:N	27:B1:87:PRO:HD3	2.20	0.56
29:B3:56:VAL:CG1	29:B3:57:GLU:N	2.65	0.56
35:BA:1763:G:H4'	35:BA:1763:G:OP1	2.05	0.56
35:BA:1899:G:O2'	35:BA:1900:A:H5''	2.05	0.56
35:BA:1998:G:O2'	35:BA:1999:C:H5'	2.06	0.56
35:BA:203:C:H2'	35:BA:204:A:C8	2.40	0.56
35:BA:2810:A:H2'	39:BE:61:ARG:HH21	1.69	0.56
35:BA:2836:U:H2'	35:BA:2837:G:C8	2.41	0.56
35:BA:292:C:N4	35:BA:348:G:H1	2.02	0.56
39:BE:3:GLY:HA3	39:BE:81:ILE:HG21	1.86	0.56
40:BF:107:LYS:O	40:BF:110:LEU:N	2.39	0.56
40:BF:139:PHE:HB3	40:BF:166:ALA:HB1	1.88	0.56
35:BA:322:A:OP2	40:BF:169:ASN:HB2	2.05	0.56
44:BN:58:ASP:C	44:BN:60:ILE:N	2.58	0.56
45:BO:62:VAL:HG11	45:BO:65:THR:HG22	1.87	0.56
46:BP:148:LEU:H	46:BP:148:LEU:HD13	1.69	0.56
52:BV:22:VAL:O	52:BV:23:GLU:CB	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:BX:21:PHE:CE1	54:BX:26:TYR:HB3	2.41	0.56
55:BY:31:LEU:HD11	55:BY:34:LYS:H	1.70	0.56
55:BY:77:PRO:O	55:BY:78:ALA:HB2	2.05	0.56
56:BZ:77:ASP:O	56:BZ:77:ASP:OD1	2.23	0.56
1:CA:1104:G:OP1	2:CB:111:ARG:HD2	2.05	0.56
1:CA:1118:C:H1'	1:CA:1179:A:C4	2.41	0.56
1:CA:1229:A:H2'	1:CA:1230:C:H6	1.71	0.56
1:CA:236:G:H2'	1:CA:237:C:C6	2.40	0.56
5:CE:133:TYR:HD1	5:CE:133:TYR:H	1.53	0.56
7:CG:121:ALA:O	7:CG:124:LEU:HB2	2.06	0.56
7:CG:148:ASN:C	7:CG:150:ALA:H	2.07	0.56
7:CG:27:ILE:HG21	7:CG:40:ALA:HB2	1.87	0.56
10:CJ:75:ILE:HG13	10:CJ:76:ASN:N	2.21	0.56
17:CQ:65:ILE:HD12	17:CQ:65:ILE:N	2.20	0.56
6:CF:89:MET:SD	18:CR:76:LEU:HD21	2.45	0.56
20:CT:63:ILE:O	20:CT:65:LYS:N	2.38	0.56
31:D5:46:CYS:SG	31:D5:48:GLU:HG3	2.46	0.56
32:D6:9:LEU:HD23	32:D6:10:LEU:N	2.20	0.56
35:DA:1184:G:O2'	35:DA:1185:C:H5'	2.05	0.56
35:DA:132:G:H5'	35:DA:132:G:C8	2.35	0.56
35:DA:1759:A:H2'	35:DA:1760:A:C8	2.41	0.56
35:DA:221:A:H4'	35:DA:222:A:O5'	2.06	0.56
35:DA:2579:C:C2'	35:DA:2580:U:H5'	2.35	0.56
35:DA:2801:A:O2'	35:DA:2895:U:H4'	2.06	0.56
35:DA:2843:G:H1	35:DA:2874:C:H42	1.54	0.56
35:DA:705:A:O2'	35:DA:706:A:H5'	2.04	0.56
35:DA:941:A:H4'	46:DP:35:HIS:CE1	2.40	0.56
36:DB:21:G:N3	36:DB:21:G:H2'	2.20	0.56
38:DD:92:ILE:HA	38:DD:107:ALA:HB2	1.87	0.56
39:DE:120:TRP:CE3	39:DE:155:LYS:HD3	2.41	0.56
43:DI:114:LEU:O	43:DI:115:ALA:HB3	2.06	0.56
43:DI:10:GLU:CD	43:DI:11:ASN:H	2.09	0.56
43:DI:37:VAL:HG13	43:DI:38:LEU:CD1	2.34	0.56
43:DI:84:GLY:O	43:DI:85:GLU:HB2	2.05	0.56
45:DO:32:TYR:HD1	45:DO:32:TYR:N	2.01	0.56
46:DP:132:LYS:O	46:DP:136:GLU:HG2	2.06	0.56
46:DP:122:PRO:HB3	46:DP:141:ALA:CB	2.36	0.56
47:DQ:81:VAL:HG23	47:DQ:82:ARG:HG2	1.86	0.56
48:DR:36:THR:HB	48:DR:40:LYS:HD2	1.86	0.56
48:DR:41:ALA:O	48:DR:43:GLU:N	2.38	0.56
53:DW:9:TYR:H	53:DW:102:HIS:CD2	2.23	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:27:VAL:HG12	55:DY:29:GLU:OE1	2.05	0.56
56:DZ:73:GLN:O	56:DZ:87:ASP:OD2	2.23	0.56
1:AA:1509:C:H2'	1:AA:1510:U:O4'	2.05	0.56
1:AA:297:G:H2'	1:AA:299:G:OP2	2.05	0.56
1:AA:359:U:H2'	1:AA:360:A:C8	2.40	0.56
1:AA:449:C:H2'	1:AA:450:G:O4'	2.05	0.56
1:AA:685:G:N2	1:AA:706:A:H61	2.03	0.56
2:AB:105:PHE:HA	2:AB:108:ILE:HG22	1.88	0.56
2:AB:69:LEU:HB3	2:AB:162:ILE:HG22	1.88	0.56
4:AD:116:GLN:NE2	4:AD:157:LEU:HD21	2.21	0.56
4:AD:163:GLU:C	4:AD:165:MET:N	2.59	0.56
4:AD:59:ARG:HH11	4:AD:59:ARG:HG2	1.70	0.56
5:AE:146:ALA:C	5:AE:148:VAL:H	2.09	0.56
5:AE:76:ILE:HG23	5:AE:77:PRO:N	2.19	0.56
11:AK:43:SER:OG	11:AK:47:VAL:HG11	2.06	0.56
12:AL:38:THR:HG22	12:AL:57:LYS:C	2.26	0.56
12:AL:86:ARG:CG	12:AL:87:GLY:H	2.18	0.56
12:AL:90:VAL:O	12:AL:92:ASP:N	2.33	0.56
12:AL:9:GLN:O	12:AL:12:ARG:N	2.39	0.56
14:AN:27:CYS:SG	14:AN:29:ARG:HB2	2.46	0.56
15:AO:50:HIS:O	15:AO:53:HIS:HB3	2.05	0.56
16:AP:71:ARG:NH1	16:AP:71:ARG:HG3	2.21	0.56
19:AS:43:GLU:O	19:AS:45:VAL:N	2.38	0.56
34:B8:38:GLY:C	34:B8:40:GLU:H	2.08	0.56
35:BA:1717:G:C3'	35:BA:1718:G:H5''	2.33	0.56
35:BA:1795:C:H1'	35:BA:1901:A:OP1	2.06	0.56
35:BA:2563:U:O2'	45:BO:28:SER:HB3	2.06	0.56
35:BA:262:A:H2'	35:BA:263:C:O4'	2.05	0.56
35:BA:2888:C:H2'	35:BA:2889:C:H6	1.69	0.56
35:BA:62:C:H2'	35:BA:63:U:H5'	1.88	0.56
35:BA:811:U:H1'	35:BA:1251:C:C5'	2.35	0.56
35:BA:853:G:H2'	35:BA:854:G:C8	2.39	0.56
41:BG:161:THR:CG2	41:BG:163:ALA:HB3	2.35	0.56
45:BO:35:VAL:H	45:BO:65:THR:HG21	1.70	0.56
50:BT:34:VAL:O	50:BT:34:VAL:HG12	2.06	0.56
56:BZ:139:VAL:O	56:BZ:141:VAL:HG12	2.04	0.56
1:CA:338:A:H2'	1:CA:339:C:H6	1.71	0.56
1:CA:597:G:H2'	1:CA:598:U:C5'	2.36	0.56
4:CD:91:SER:HA	4:CD:94:LEU:HD12	1.87	0.56
5:CE:19:MET:O	5:CE:20:GLN:HB2	2.04	0.56
7:CG:29:LYS:CB	7:CG:105:VAL:HG21	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:63:LEU:HB3	8:CH:65:TYR:CE1	2.40	0.56
1:CA:599:C:O3'	8:CH:96:GLY:HA2	2.06	0.56
10:CJ:48:THR:HG23	10:CJ:62:HIS:N	2.21	0.56
19:CS:53:ASN:ND2	19:CS:53:ASN:N	2.54	0.56
20:CT:63:ILE:HD12	20:CT:81:LYS:CG	2.35	0.56
22:CV:30:A:C4	22:CV:31:U:C5	2.91	0.56
23:CW:28:U:O5'	23:CW:28:U:H6	1.88	0.56
25:CY:111:ARG:O	25:CY:115:VAL:HG23	2.05	0.56
33:D7:5:TRP:NE1	33:D7:7:PRO:HG3	2.20	0.56
34:D8:23:VAL:CG1	34:D8:46:ARG:NH1	2.69	0.56
35:DA:173:G:H2'	35:DA:174:C:C6	2.41	0.56
35:DA:1799:G:H4'	35:DA:1800:C:O5'	2.05	0.56
35:DA:2286:A:H4'	35:DA:2287:A:O4'	2.06	0.56
35:DA:237:C:H2'	35:DA:238:C:C6	2.40	0.56
35:DA:271(L):U:H4'	35:DA:271(M):G:C5	2.40	0.56
35:DA:302:C:H42	35:DA:315:G:H1	1.54	0.56
35:DA:444:C:O5'	51:DU:2:PRO:HD3	2.05	0.56
37:DC:56:GLN:HE22	37:DC:169:GLY:H	1.53	0.56
38:DD:186:HIS:HB3	38:DD:189:CYS:SG	2.45	0.56
38:DD:97:TYR:HB2	38:DD:101:GLU:O	2.06	0.56
39:DE:113:PHE:CE2	39:DE:158:GLY:HA2	2.41	0.56
39:DE:63:LEU:O	39:DE:64:LYS:C	2.44	0.56
39:DE:78:LEU:CD2	39:DE:78:LEU:N	2.69	0.56
40:DF:183:VAL:O	40:DF:187:VAL:HG23	2.06	0.56
41:DG:105:LYS:HB3	41:DG:142:PRO:HG3	1.88	0.56
42:DH:159:GLU:HA	42:DH:159:GLU:OE1	2.06	0.56
44:DN:67:LEU:C	44:DN:69:GLN:H	2.09	0.56
46:DP:30:THR:O	46:DP:32:THR:N	2.39	0.56
48:DR:45:ARG:HG3	48:DR:46:GLY:N	2.10	0.56
49:DS:66:ALA:HA	49:DS:69:VAL:HG12	1.87	0.56
52:DV:34:GLU:CB	52:DV:62:LEU:HD12	2.36	0.56
52:DV:58:VAL:HG12	52:DV:101:GLY:O	2.06	0.56
53:DW:29:LEU:HD23	53:DW:30:GLU:N	2.21	0.56
53:DW:87:PRO:HA	53:DW:93:ALA:CB	2.35	0.56
54:DX:51:VAL:CG1	54:DX:80:ILE:H	2.18	0.56
55:DY:74:PRO:O	55:DY:75:ILE:CB	2.54	0.56
56:DZ:8:TYR:H	56:DZ:62:PRO:CD	2.18	0.56
1:AA:1223:C:OP1	1:AA:1224:G:H2'	2.06	0.56
1:AA:1296:C:H5'	1:AA:1297:C:OP2	2.06	0.56
1:AA:1320:C:H5'	19:AS:70:LYS:HE3	1.88	0.56
1:AA:1325:C:H2'	1:AA:1326:C:H5'	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1367:C:OP1	9:AI:115:GLY:N	2.39	0.56
1:AA:1516:G:H2'	1:AA:1518:A:OP2	2.06	0.56
1:AA:522:C:H2'	1:AA:523:A:C8	2.41	0.56
1:AA:579:G:H5'	1:AA:728:A:H1'	1.87	0.56
5:AE:19:MET:O	5:AE:20:GLN:HB2	2.05	0.56
6:AF:27:GLN:HE21	6:AF:27:GLN:HA	1.69	0.56
7:AG:36:LYS:HA	7:AG:39:ALA:CB	2.36	0.56
7:AG:26:PHE:HB2	7:AG:62:PHE:CZ	2.41	0.56
19:AS:4:SER:N	19:AS:6:LYS:HZ1	2.03	0.56
20:AT:72:LEU:HB3	20:AT:76:ALA:HB1	1.88	0.56
25:AY:101:ILE:N	25:AY:101:ILE:HD12	2.20	0.56
28:B2:53:LEU:O	28:B2:54:LYS:CB	2.52	0.56
29:B3:43:ILE:O	29:B3:47:VAL:HG23	2.06	0.56
33:B7:34:ARG:O	33:B7:36:GLN:N	2.38	0.56
35:BA:1189:A:C2	35:BA:1190:G:H1'	2.40	0.56
35:BA:1197:G:H5'	35:BA:1227:G:O2'	2.06	0.56
35:BA:121:G:H2'	35:BA:122:G:H8	1.71	0.56
35:BA:1322:A:O2'	35:BA:1323:U:H5'	2.05	0.56
35:BA:1553:A:H2'	35:BA:1554:A:H5''	1.88	0.56
35:BA:1688:U:H5'	35:BA:1689:A:OP1	2.06	0.56
35:BA:2308:G:O6	35:BA:2310:A:H2'	2.05	0.56
35:BA:2328:A:H2'	35:BA:2329:G:C8	2.41	0.56
35:BA:27:G:N2	35:BA:512:G:C2'	2.69	0.56
35:BA:28:A:H1'	35:BA:513:A:C2	2.41	0.56
35:BA:543:C:H6	35:BA:547:A:C8	2.24	0.56
35:BA:658:C:H2'	35:BA:659:C:H6	1.70	0.56
35:BA:66:C:C2'	35:BA:67:U:H5'	2.35	0.56
35:BA:684:G:H22	35:BA:787:U:H2'	1.70	0.56
35:BA:902:C:H2'	35:BA:903:C:C6	2.41	0.56
38:BD:211:ARG:HA	38:BD:214:TRP:CD2	2.40	0.56
39:BE:52:LEU:HD12	39:BE:53:PRO:CD	2.36	0.56
42:BH:85:LYS:HE3	42:BH:144:VAL:HB	1.88	0.56
43:BI:41:GLU:HA	43:BI:44:LEU:HB3	1.88	0.56
44:BN:3:THR:O	44:BN:5:VAL:N	2.28	0.56
45:BO:31:LYS:HB3	45:BO:32:TYR:CD1	2.41	0.56
48:BR:104:ARG:O	48:BR:106:GLY:N	2.38	0.56
48:BR:24:GLN:CB	48:BR:44:LEU:HD21	2.32	0.56
49:BS:66:ALA:HA	49:BS:69:VAL:HG12	1.88	0.56
50:BT:94:ALA:CB	50:BT:99:LEU:HD23	2.35	0.56
51:BU:65:ILE:HD12	51:BU:65:ILE:H	1.70	0.56
1:CA:1131:G:H2'	1:CA:1132:C:C6	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1293:G:HO2'	1:CA:1294:G:H8	1.52	0.56
1:CA:436:C:O2'	1:CA:437:U:P	2.64	0.56
1:CA:501:C:O2'	1:CA:502:G:H5'	2.05	0.56
1:CA:64:G:H3'	1:CA:64:G:OP1	2.05	0.56
2:CB:169:LYS:HB3	2:CB:170:GLU:OE2	2.06	0.56
2:CB:81:VAL:HG22	2:CB:215:LEU:HG	1.88	0.56
3:CC:16:ARG:NH1	3:CC:16:ARG:CB	2.68	0.56
4:CD:108:LEU:O	4:CD:110:PHE:CD1	2.59	0.56
5:CE:147:ASP:HA	5:CE:150:ARG:NH1	2.16	0.56
6:CF:23:LYS:O	6:CF:27:GLN:HG2	2.05	0.56
8:CH:109:ILE:CG1	8:CH:110:ALA:N	2.69	0.56
12:CL:38:THR:HG22	12:CL:57:LYS:C	2.25	0.56
25:CY:10:THR:C	25:CY:12:SER:N	2.59	0.56
28:D2:26:ARG:HD3	54:DX:5:TYR:HB3	1.88	0.56
35:DA:2020:A:N1	35:DA:2034:U:O4	2.38	0.56
35:DA:212:G:O2'	35:DA:213:A:H5'	2.05	0.56
35:DA:2589:A:H2'	35:DA:2590:A:C8	2.41	0.56
35:DA:2682:U:O2	39:DE:22:PRO:HB3	2.06	0.56
35:DA:2687:U:O2'	35:DA:2688:U:H5'	2.05	0.56
35:DA:2692:C:O2'	35:DA:2693:A:H5'	2.06	0.56
35:DA:272:G:H1'	35:DA:272(B):G:O4'	2.06	0.56
35:DA:568:U:H2'	35:DA:570:G:OP2	2.06	0.56
35:DA:61:G:O2'	35:DA:62:C:H5'	2.05	0.56
38:DD:79:VAL:HG11	38:DD:112:GLN:O	2.06	0.56
35:DA:1843:C:C1'	38:DD:255:LYS:HZ3	2.18	0.56
38:DD:265:PRO:HG2	38:DD:266:SER:N	2.17	0.56
40:DF:125:LEU:HB3	40:DF:196:LEU:HD21	1.88	0.56
40:DF:184:TYR:CD2	40:DF:185:ASP:N	2.73	0.56
41:DG:132:ASN:ND2	41:DG:133:LEU:H	2.04	0.56
41:DG:51:ARG:HH12	41:DG:53:LEU:HG	1.61	0.56
43:DI:6:LEU:O	43:DI:15:VAL:HG12	2.06	0.56
45:DO:87:ILE:HD12	45:DO:92:GLU:N	2.20	0.56
46:DP:16:ARG:NE	46:DP:18:ARG:HB2	2.21	0.56
52:DV:18:LEU:O	52:DV:19:LYS:O	2.24	0.56
55:DY:83:THR:HG22	55:DY:84:ARG:N	2.21	0.56
1:AA:1226:C:H4'	1:AA:1227:A:OP1	2.06	0.56
1:AA:1292:U:H2'	1:AA:1293:G:H8	1.69	0.56
1:AA:189(H):G:H2'	1:AA:189(I):G:C8	2.41	0.56
1:AA:437:U:O2'	1:AA:438:G:H5'	2.06	0.56
1:AA:539:A:H2'	1:AA:540:G:H8	1.70	0.56
1:AA:802:A:H2'	1:AA:803:G:H5'	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:899:C:H2'	1:AA:900:A:O4'	2.05	0.56
1:AA:939:G:H5''	7:AG:102:ARG:HH12	1.71	0.56
4:AD:58:LEU:O	4:AD:59:ARG:C	2.43	0.56
6:AF:70:ASP:CG	6:AF:71:ARG:H	2.09	0.56
6:AF:80:ARG:CG	6:AF:88:VAL:HB	2.36	0.56
7:AG:83:ALA:C	7:AG:84:ASN:HD22	2.08	0.56
13:AM:40:ASN:ND2	13:AM:43:THR:HG23	2.21	0.56
15:AO:85:LEU:HD12	15:AO:87:ILE:HD11	1.88	0.56
16:AP:71:ARG:HH11	16:AP:71:ARG:HG3	1.69	0.56
17:AQ:12:SER:HB3	17:AQ:20:THR:OG1	2.06	0.56
28:B2:15:LYS:O	28:B2:19:VAL:HG23	2.06	0.56
34:B8:30:ARG:O	34:B8:30:ARG:HG3	2.05	0.56
35:BA:1021:A:H8	35:BA:1021:A:H3'	1.70	0.56
35:BA:2128:C:O2'	35:BA:2163:C:OP1	2.18	0.56
35:BA:2687:U:O2'	35:BA:2688:U:H5'	2.05	0.56
35:BA:2713:A:C3'	35:BA:2714:G:H5'	2.36	0.56
35:BA:483:A:H3'	35:BA:484:C:H6	1.71	0.56
35:BA:52:A:O2'	35:BA:53:A:H5'	2.05	0.56
36:BB:74:U:H2'	36:BB:75:G:C5'	2.18	0.56
38:BD:145:VAL:HB	38:BD:155:LEU:HB2	1.88	0.56
40:BF:110:LEU:O	40:BF:113:ALA:HB3	2.05	0.56
41:BG:140:ILE:HD11	41:BG:141:PHE:CE2	2.41	0.56
41:BG:91:ARG:CG	41:BG:92:VAL:N	2.69	0.56
45:BO:4:PRO:O	45:BO:5:GLN:HB3	2.04	0.56
45:BO:61:VAL:O	45:BO:84:ALA:CB	2.47	0.56
46:BP:107:LYS:O	46:BP:109:GLY:N	2.38	0.56
46:BP:30:THR:O	46:BP:32:THR:N	2.39	0.56
35:BA:2009:G:H1'	48:BR:107:ASP:O	2.06	0.56
51:BU:52:ARG:O	51:BU:54:LYS:N	2.38	0.56
52:BV:18:LEU:O	52:BV:19:LYS:O	2.23	0.56
52:BV:32:THR:CG2	52:BV:33:VAL:H	2.14	0.56
52:BV:40:LEU:HD13	52:BV:40:LEU:O	2.06	0.56
53:BW:74:ALA:C	53:BW:75:TYR:CD1	2.79	0.56
56:BZ:53:ILE:HG22	56:BZ:71:VAL:HB	1.88	0.56
1:CA:1208:C:H2'	1:CA:1209:C:C6	2.41	0.56
1:CA:1223:C:OP1	1:CA:1224:G:H2'	2.06	0.56
1:CA:1276:G:H2'	1:CA:1277:C:H5'	1.87	0.56
1:CA:284:G:H2'	1:CA:285:G:C8	2.41	0.56
2:CB:159:PRO:O	2:CB:161:ALA:N	2.38	0.56
2:CB:238:LEU:O	2:CB:239:VAL:C	2.43	0.56
4:CD:25:ARG:C	4:CD:27:TYR:H	2.09	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:50:GLU:OE2	5:CE:51:VAL:HG23	2.05	0.56
11:CK:23:ALA:HA	11:CK:28:THR:CG2	2.36	0.56
6:CF:100:ASN:HB3	18:CR:28:GLU:HA	1.88	0.56
19:CS:10:PHE:HE2	19:CS:70:LYS:HZ1	1.53	0.56
20:CT:14:LYS:HA	20:CT:17:ARG:HE	1.71	0.56
22:CV:36:A:H3'	22:CV:37:A:H8	1.71	0.56
27:D1:9:GLY:N	27:D1:48:LYS:HZ2	2.03	0.56
28:D2:30:ARG:HH11	28:D2:30:ARG:HG3	1.70	0.56
28:D2:56:GLN:HG3	35:DA:76:C:O2'	2.06	0.56
33:D7:16:HIS:HD1	33:D7:21:ARG:HH22	1.52	0.56
35:DA:1009:A:H2'	35:DA:1010:A:C8	2.40	0.56
35:DA:1417:C:O2'	35:DA:1418:G:H5'	2.05	0.56
35:DA:1485:G:H2'	35:DA:1486:A:H8	1.71	0.56
35:DA:1688:U:H1'	35:DA:1701:A:C6	2.41	0.56
35:DA:2307:G:H21	35:DA:2308:G:H5'	1.70	0.56
38:DD:92:ILE:HA	38:DD:107:ALA:CB	2.36	0.56
38:DD:125:ILE:CD1	38:DD:125:ILE:H	2.19	0.56
38:DD:257:LEU:C	38:DD:257:LEU:CD2	2.74	0.56
38:DD:270:ILE:O	38:DD:270:ILE:HD12	2.05	0.56
35:DA:1490:A:C2	38:DD:75:ILE:HD12	2.40	0.56
39:DE:120:TRP:NE1	39:DE:155:LYS:HB3	2.21	0.56
44:DN:56:ASN:HA	44:DN:125:GLY:N	2.20	0.56
35:DA:1030:G:OP2	47:DQ:128:LYS:HE2	2.06	0.56
47:DQ:9:TYR:CG	47:DQ:9:TYR:O	2.58	0.56
50:DT:99:LEU:O	50:DT:99:LEU:HD13	2.06	0.56
53:DW:59:VAL:HG12	53:DW:60:ASN:N	2.20	0.56
54:DX:9:LEU:HG	54:DX:29:TRP:O	2.06	0.56
55:DY:31:LEU:HG	55:DY:34:LYS:HB2	1.88	0.56
56:DZ:59:LEU:H	56:DZ:59:LEU:HD23	1.71	0.56
1:AA:652:U:H1'	1:AA:653:A:H2	1.69	0.56
2:AB:222:ILE:HG23	2:AB:223:ILE:N	2.20	0.56
3:AC:182:ILE:HG23	3:AC:203:PHE:CA	2.36	0.56
4:AD:22:LYS:HG3	4:AD:26:CYS:SG	2.46	0.56
4:AD:80:GLU:C	4:AD:84:LYS:HZ3	2.09	0.56
8:AH:11:THR:HG23	8:AH:14:ARG:NH1	2.15	0.56
9:AI:16:ARG:O	9:AI:63:ILE:HG22	2.06	0.56
1:AA:552:U:H4'	12:AL:86:ARG:HG2	1.88	0.56
16:AP:14:ASN:H	16:AP:15:PRO:HD3	1.67	0.56
19:AS:15:LEU:O	19:AS:19:VAL:HG23	2.06	0.56
19:AS:36:ARG:NH1	19:AS:75:ALA:HB3	2.20	0.56
28:B2:21:LEU:HD13	28:B2:50:ILE:HG22	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2092:U:C4'	35:BA:2093:G:H5''	2.31	0.56
35:BA:648:G:H4'	35:BA:2351:G:H5''	1.87	0.56
35:BA:2589:A:H2'	35:BA:2590:A:H8	1.71	0.56
35:BA:2675:A:OP1	45:BO:31:LYS:HB2	2.05	0.56
35:BA:993:G:OP1	51:BU:50:ARG:NH2	2.39	0.56
39:BE:12:THR:CG2	50:BT:8:LYS:HE2	2.36	0.56
39:BE:88:GLY:O	39:BE:89:ASP:HB2	2.06	0.56
41:BG:101:ILE:HG12	41:BG:105:LYS:HE3	1.88	0.56
41:BG:160:VAL:HG12	41:BG:161:THR:N	2.21	0.56
41:BG:43:LEU:HD12	41:BG:153:ARG:HD3	1.87	0.56
43:BI:109:ILE:CD1	43:BI:111:PRO:HD3	2.35	0.56
43:BI:115:ALA:O	43:BI:128:LEU:HD23	2.06	0.56
44:BN:10:GLU:HG3	44:BN:11:PRO:HD2	1.86	0.56
44:BN:137:LYS:HG2	44:BN:138:LEU:N	2.20	0.56
50:BT:28:VAL:CG1	50:BT:46:GLU:HA	2.35	0.56
51:BU:92:ARG:CG	51:BU:94:ASN:HB3	2.36	0.56
52:BV:52:VAL:O	52:BV:54:GLY:N	2.38	0.56
35:BA:993:G:H5'	52:BV:75:PHE:HZ	1.71	0.56
52:BV:78:LYS:HD3	52:BV:78:LYS:C	2.27	0.56
1:CA:1109:C:H2'	1:CA:1110:A:O4'	2.06	0.56
1:CA:1415:G:C4	1:CA:1486:G:C2	2.93	0.56
1:CA:553:A:H2'	1:CA:554:C:H6	1.68	0.56
1:CA:726:C:H2'	1:CA:727:G:C8	2.40	0.56
1:CA:802:A:H2'	1:CA:803:G:H5'	1.88	0.56
2:CB:36:ARG:N	2:CB:41:ILE:HD13	2.21	0.56
2:CB:71:VAL:O	2:CB:164:VAL:HG22	2.06	0.56
3:CC:66:VAL:O	3:CC:66:VAL:HG12	2.06	0.56
4:CD:104:VAL:O	4:CD:108:LEU:HD13	2.06	0.56
4:CD:18:LYS:CE	4:CD:31:CYS:HB3	2.35	0.56
5:CE:150:ARG:NH1	5:CE:150:ARG:CB	2.69	0.56
6:CF:37:VAL:HA	6:CF:65:VAL:CG1	2.36	0.56
7:CG:83:ALA:C	7:CG:84:ASN:HD22	2.09	0.56
13:CM:40:ASN:ND2	13:CM:43:THR:HG23	2.21	0.56
15:CO:53:HIS:HE1	15:CO:57:LEU:HD21	1.71	0.56
16:CP:8:ARG:NH2	16:CP:15:PRO:HG3	2.21	0.56
16:CP:45:THR:HG22	16:CP:47:ASP:N	2.19	0.56
25:CY:30:THR:CG2	25:CY:179:LYS:HE3	2.36	0.56
26:D0:21:LEU:CD1	26:D0:41:ARG:HD3	2.36	0.56
32:D6:20:ASN:ND2	32:D6:21:TYR:N	2.52	0.56
32:D6:12:GLU:HB3	32:D6:23:THR:HG22	1.87	0.56
33:D7:19:ARG:HG2	33:D7:19:ARG:HH11	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1197:G:H2'	35:DA:1198:U:H6	1.69	0.56
35:DA:1447:G:H2'	35:DA:1448:G:H8	1.71	0.56
35:DA:1744:C:O2'	35:DA:1745:C:H5'	2.06	0.56
35:DA:1930:G:H22	35:DA:1968:G:C2'	2.19	0.56
35:DA:2081:C:O2'	35:DA:2082:A:H5'	2.06	0.56
35:DA:2221:G:H5'	35:DA:2222:G:OP2	2.06	0.56
35:DA:234:C:H2'	35:DA:235:U:C6	2.41	0.56
35:DA:2870:C:C2'	35:DA:2871:C:H5'	2.35	0.56
35:DA:627:A:H4'	35:DA:628:G:OP1	2.05	0.56
38:DD:135:PHE:CD1	38:DD:135:PHE:N	2.71	0.56
38:DD:257:LEU:HD23	38:DD:258:LYS:N	2.21	0.56
39:DE:75:VAL:C	39:DE:77:ILE:H	2.08	0.56
41:DG:129:GLY:C	41:DG:131:TYR:N	2.58	0.56
42:DH:105:LEU:CD2	42:DH:113:VAL:HB	2.36	0.56
42:DH:38:SER:O	42:DH:40:GLU:N	2.39	0.56
42:DH:41:MET:HE3	42:DH:54:ARG:HA	1.88	0.56
45:DO:62:VAL:HG12	45:DO:63:VAL:N	2.21	0.56
51:DU:49:HIS:HA	51:DU:52:ARG:HB2	1.87	0.56
53:DW:35:ILE:HG22	53:DW:36:LEU:N	2.20	0.56
54:DX:65:ARG:NE	54:DX:66:LEU:H	2.03	0.56
55:DY:6:HIS:N	55:DY:6:HIS:ND1	2.53	0.56
47:DQ:141:GLN:CD	56:DZ:89:PHE:HB3	2.27	0.56
1:AA:1261:A:H62	1:AA:1274:G:H21	1.54	0.55
1:AA:1427:U:H2'	1:AA:1428:A:C8	2.41	0.55
1:AA:498:U:C2'	1:AA:498:U:O2	2.55	0.55
2:AB:169:LYS:HB3	2:AB:170:GLU:OE2	2.07	0.55
1:AA:18:C:P	5:AE:127:ASN:HD21	2.30	0.55
5:AE:139:LEU:HA	5:AE:142:LEU:CD1	2.33	0.55
7:AG:79:ARG:NE	7:AG:84:ASN:ND2	2.52	0.55
8:AH:109:ILE:CG1	8:AH:110:ALA:N	2.67	0.55
1:AA:691:G:H1	11:AK:52:GLY:HA2	1.71	0.55
26:B0:56:ASP:CG	26:B0:58:THR:HG1	2.09	0.55
27:B1:58:ILE:CD1	27:B1:87:PRO:HB3	2.36	0.55
29:B3:4:LEU:HD21	29:B3:56:VAL:HG13	1.87	0.55
35:BA:1002:G:H2'	35:BA:1003:G:O4'	2.05	0.55
35:BA:1423:G:H2'	35:BA:1424:G:H8	1.70	0.55
35:BA:154:G:H2'	35:BA:154(A):C:O2	2.06	0.55
35:BA:1824:G:H2'	35:BA:1825:A:H8	1.71	0.55
35:BA:1960:A:O2'	35:BA:1961:C:H5'	2.07	0.55
35:BA:2111:C:H1'	35:BA:2118:U:O4'	2.06	0.55
35:BA:302:C:O2'	35:BA:303:U:H5'	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:586:A:H2	35:BA:809:G:N3	2.04	0.55
35:BA:94:C:H5'	35:BA:94(A):G:OP2	2.07	0.55
38:BD:143:HIS:HB3	38:BD:194:GLY:O	2.05	0.55
35:BA:778:G:C5'	38:BD:48:ARG:HD2	2.36	0.55
38:BD:64:ILE:HG23	38:BD:64:ILE:O	2.05	0.55
40:BF:60:SER:OG	40:BF:61:GLY:N	2.37	0.55
41:BG:166:ASP:O	41:BG:170:ARG:HB2	2.06	0.55
41:BG:69:ALA:HB1	41:BG:91:ARG:HE	1.71	0.55
42:BH:153:LYS:N	42:BH:153:LYS:HD3	2.22	0.55
46:BP:13:ASN:HD22	46:BP:13:ASN:N	2.02	0.55
47:BQ:9:TYR:CG	47:BQ:9:TYR:O	2.59	0.55
51:BU:91:ASP:O	51:BU:95:LEU:HB2	2.07	0.55
52:BV:71:LEU:O	52:BV:90:PRO:HA	2.06	0.55
1:CA:112:G:O2'	1:CA:113:G:H5'	2.05	0.55
1:CA:1226:C:H4'	1:CA:1227:A:OP1	2.06	0.55
1:CA:1244:C:H2'	1:CA:1245:A:H8	1.71	0.55
1:CA:1469:G:H2'	1:CA:1470:G:H8	1.71	0.55
1:CA:965:A:C2	1:CA:969:A:C2	2.94	0.55
2:CB:101:MET:O	2:CB:108:ILE:HG21	2.05	0.55
3:CC:153:VAL:HG12	3:CC:154:SER:N	2.21	0.55
6:CF:71:ARG:O	6:CF:73:ASN:N	2.38	0.55
7:CG:36:LYS:HA	7:CG:39:ALA:CB	2.36	0.55
7:CG:79:ARG:NE	7:CG:84:ASN:ND2	2.51	0.55
12:CL:89:ARG:O	12:CL:89:ARG:HD3	2.06	0.55
16:CP:60:LEU:C	16:CP:62:VAL:H	2.09	0.55
17:CQ:29:HIS:HE1	17:CQ:31:LEU:HB3	1.71	0.55
23:CW:22:A:N6	23:CW:47:G:H1'	2.21	0.55
27:D1:21:ARG:HD3	27:D1:22:GLY:N	2.20	0.55
27:D1:71:TYR:HA	27:D1:74:VAL:HG23	1.87	0.55
28:D2:60:LEU:HB3	28:D2:61:LEU:HD23	1.87	0.55
35:DA:1158:C:HO2'	35:DA:1159:U:H6	1.53	0.55
35:DA:1189:A:C2	35:DA:1190:G:H1'	2.41	0.55
35:DA:1215:G:H2'	35:DA:1216:G:C8	2.41	0.55
35:DA:147:U:H2'	35:DA:148:C:H6	1.71	0.55
35:DA:2233:U:H2'	35:DA:2234:G:C8	2.41	0.55
35:DA:2266:A:H4'	35:DA:2267:A:C2	2.41	0.55
35:DA:2380:C:H2'	35:DA:2381:C:C6	2.41	0.55
35:DA:2476:A:N3	35:DA:2477:C:H5'	2.21	0.55
27:D1:66:HIS:NE2	35:DA:372:G:H3'	2.21	0.55
35:DA:674:G:P	40:DF:54:ARG:HH22	2.29	0.55
41:DG:59:GLU:O	41:DG:63:ILE:HG23	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DI:87:LYS:HZ3	43:DI:121:LYS:HG2	1.70	0.55
47:DQ:140:ALA:HB3	56:DZ:53:ILE:HD12	1.87	0.55
48:DR:70:LEU:O	48:DR:71:GLN:HB2	2.06	0.55
50:DT:102:ILE:CB	50:DT:110:ILE:HD11	2.35	0.55
51:DU:91:ASP:OD2	51:DU:96:ALA:HA	2.05	0.55
52:DV:22:VAL:O	52:DV:23:GLU:HB2	2.04	0.55
35:DA:993:G:H5'	52:DV:75:PHE:HZ	1.68	0.55
54:DX:57:LEU:HB2	54:DX:76:ARG:CD	2.33	0.55
54:DX:82:GLN:HB3	54:DX:85:PRO:HG2	1.88	0.55
56:DZ:57:ILE:HG22	56:DZ:59:LEU:HD23	1.87	0.55
56:DZ:24:LEU:CD1	56:DZ:86:VAL:HG23	2.30	0.55
1:AA:1365:G:O2'	1:AA:1366:C:H5'	2.07	0.55
1:AA:1514:C:H2'	1:AA:1515:C:H6	1.69	0.55
1:AA:192:U:H2'	1:AA:193:C:C6	2.41	0.55
1:AA:553:A:H2'	1:AA:554:C:H6	1.71	0.55
1:AA:597:G:H2'	1:AA:598:U:C5'	2.36	0.55
2:AB:159:PRO:O	2:AB:161:ALA:N	2.39	0.55
2:AB:22:LYS:HZ3	2:AB:22:LYS:HA	1.70	0.55
3:AC:87:LEU:HB3	3:AC:101:LEU:HD11	1.88	0.55
8:AH:9:MET:O	8:AH:13:ILE:HG12	2.05	0.55
8:AH:91:ARG:HH11	8:AH:91:ARG:CG	2.19	0.55
9:AI:117:HIS:HB2	9:AI:121:ARG:HD2	1.88	0.55
10:AJ:62:HIS:N	10:AJ:62:HIS:CD2	2.74	0.55
13:AM:77:ASN:O	13:AM:81:LEU:HG	2.05	0.55
17:AQ:58:GLU:HB2	17:AQ:74:LEU:HB3	1.88	0.55
18:AR:36:ASN:HB2	18:AR:39:VAL:HG21	1.88	0.55
25:AY:15:GLN:HA	25:AY:168:PHE:CE2	2.41	0.55
25:AY:93:SER:O	25:AY:99:LEU:HD23	2.06	0.55
27:B1:45:ASN:O	27:B1:46:LEU:C	2.44	0.55
28:B2:14:ARG:O	28:B2:16:LEU:N	2.39	0.55
35:BA:149:A:H2'	35:BA:150:C:C6	2.41	0.55
35:BA:1759:A:H2'	35:BA:1760:A:C8	2.40	0.55
35:BA:2246:G:H2'	35:BA:2247:A:H8	1.72	0.55
35:BA:2528:U:H2'	35:BA:2530:A:O5'	2.05	0.55
35:BA:302:C:H42	35:BA:315:G:H1	1.53	0.55
35:BA:481:G:HO2'	35:BA:482:A:P	2.29	0.55
38:BD:96:HIS:HA	38:BD:102:LYS:HG2	1.87	0.55
39:BE:11:MET:HB2	39:BE:23:VAL:O	2.06	0.55
39:BE:47:VAL:HG12	39:BE:49:LEU:HD21	1.88	0.55
40:BF:154:VAL:HB	40:BF:173:VAL:HG22	1.88	0.55
40:BF:57:VAL:HG12	40:BF:58:ALA:N	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:39:ILE:HD11	41:BG:155:MET:SD	2.46	0.55
44:BN:67:LEU:C	44:BN:69:GLN:H	2.10	0.55
48:BR:28:LEU:HA	48:BR:34:ILE:HD11	1.87	0.55
50:BT:13:ARG:HH12	50:BT:15:VAL:HG12	1.67	0.55
54:BX:52:VAL:H	54:BX:80:ILE:CG2	2.20	0.55
56:BZ:61:LEU:CB	56:BZ:65:GLN:HB3	2.35	0.55
1:CA:1347:G:C2	1:CA:1373:G:H2'	2.41	0.55
1:CA:180:U:H2'	1:CA:181:G:H5''	1.87	0.55
1:CA:363:A:C2	12:CL:31:PRO:HG2	2.42	0.55
1:CA:674:G:H2'	1:CA:675:A:C8	2.35	0.55
1:CA:741:G:H5'	15:CO:39:LEU:CD2	2.36	0.55
1:CA:859:A:H2'	1:CA:860:A:O4'	2.05	0.55
1:CA:959:A:H2'	1:CA:960:U:C4'	2.36	0.55
4:CD:128:VAL:C	4:CD:130:GLY:N	2.59	0.55
5:CE:10:MET:HG2	5:CE:10:MET:O	2.05	0.55
5:CE:131:ILE:N	5:CE:131:ILE:HD13	2.21	0.55
6:CF:39:LYS:HG2	6:CF:40:VAL:N	2.21	0.55
1:CA:939:G:H5''	7:CG:102:ARG:CZ	2.37	0.55
11:CK:17:GLY:HA3	11:CK:79:SER:O	2.07	0.55
13:CM:91:ARG:HG3	13:CM:98:VAL:HG13	1.88	0.55
15:CO:33:THR:OG1	15:CO:63:ARG:HD2	2.07	0.55
18:CR:58:LEU:N	18:CR:58:LEU:HD12	2.13	0.55
20:CT:8:ARG:HD2	20:CT:8:ARG:N	2.21	0.55
25:CY:156:ARG:O	25:CY:157:ALA:C	2.44	0.55
25:CY:35:PRO:HA	25:CY:66:LEU:CD2	2.36	0.55
27:D1:19:GLN:HE21	35:DA:379:G:N2	2.03	0.55
27:D1:76:ARG:O	27:D1:77:ALA:CB	2.54	0.55
34:D8:3:LYS:O	34:D8:4:MET:O	2.23	0.55
35:DA:127:A:H5''	35:DA:128:C:O4'	2.06	0.55
35:DA:1776:G:C2	35:DA:1777:U:C6	2.94	0.55
35:DA:1778:U:H2'	35:DA:1779:U:C6	2.41	0.55
35:DA:1821:A:H2'	35:DA:1822:G:H8	1.71	0.55
35:DA:1839:G:H8	35:DA:1839:G:H5'	1.70	0.55
35:DA:1848:A:H2'	35:DA:1849:G:H8	1.71	0.55
35:DA:13:A:C2	35:DA:526:A:C5	2.94	0.55
35:DA:860:U:O2'	35:DA:861:A:H5'	2.05	0.55
37:DC:99:ILE:O	37:DC:99:ILE:HG22	2.06	0.55
38:DD:264:LYS:HE3	38:DD:266:SER:HB2	1.88	0.55
39:DE:141:ILE:HG12	39:DE:142:GLY:H	1.69	0.55
39:DE:81:ILE:O	39:DE:82:ARG:O	2.24	0.55
41:DG:102:PHE:HA	41:DG:105:LYS:NZ	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:63:VAL:HB	45:DO:102:VAL:HG12	1.88	0.55
48:DR:28:LEU:HA	48:DR:34:ILE:HD11	1.88	0.55
50:DT:27:THR:OG1	50:DT:28:VAL:N	2.36	0.55
50:DT:28:VAL:O	50:DT:29:ARG:HD3	2.07	0.55
52:DV:33:VAL:HG13	52:DV:62:LEU:H	1.70	0.55
53:DW:8:ARG:HG3	53:DW:8:ARG:HH11	1.71	0.55
54:DX:14:SER:O	54:DX:17:ALA:N	2.40	0.55
54:DX:35:THR:O	54:DX:39:ILE:HG23	2.06	0.55
47:DQ:137:TYR:HE2	56:DZ:76:LEU:HD21	1.70	0.55
1:AA:1508:G:H2'	1:AA:1509:C:C6	2.42	0.55
4:AD:190:ASP:O	4:AD:191:ARG:C	2.44	0.55
5:AE:43:LEU:HD12	5:AE:44:GLY:N	2.21	0.55
7:AG:29:LYS:CB	7:AG:105:VAL:HG21	2.35	0.55
15:AO:67:LEU:HD22	15:AO:78:TYR:HE1	1.72	0.55
16:AP:20:VAL:CG2	16:AP:21:VAL:N	2.68	0.55
16:AP:48:TRP:HE3	16:AP:49:LEU:N	2.03	0.55
23:AW:73:A:H5'	23:AW:74:A:P	2.47	0.55
35:BA:1386:C:O2'	35:BA:1387:C:H5'	2.07	0.55
35:BA:1790:C:H5''	35:BA:1791:A:OP1	2.05	0.55
35:BA:2364:C:O2'	35:BA:2365:G:H5'	2.07	0.55
35:BA:2467:C:O2'	35:BA:2468:G:H5'	2.06	0.55
35:BA:523:C:C2'	35:BA:524:U:H5'	2.36	0.55
35:BA:565:C:O3'	52:BV:81:TYR:CE1	2.60	0.55
35:BA:763:G:C4	35:BA:765:G:C8	2.94	0.55
35:BA:827:U:H2'	35:BA:2068:U:C2	2.42	0.55
39:BE:77:ILE:HG22	39:BE:79:ARG:HD2	1.88	0.55
41:BG:138:GLN:OE1	41:BG:153:ARG:HB2	2.06	0.55
41:BG:47:LYS:O	41:BG:51:ARG:HG2	2.06	0.55
42:BH:55:PRO:HG2	42:BH:61:HIS:NE2	2.22	0.55
43:BI:58:LEU:C	43:BI:60:GLU:H	2.10	0.55
45:BO:13:ASN:HD22	45:BO:97:ARG:HB2	1.71	0.55
46:BP:96:THR:O	46:BP:99:LEU:HB3	2.06	0.55
49:BS:93:LYS:HD2	49:BS:93:LYS:O	2.06	0.55
51:BU:21:ALA:CB	51:BU:35:ALA:HB1	2.36	0.55
55:BY:31:LEU:HG	55:BY:34:LYS:HB2	1.88	0.55
56:BZ:127:LYS:N	56:BZ:164:ALA:CB	2.66	0.55
47:BQ:141:GLN:CD	56:BZ:89:PHE:HB3	2.27	0.55
1:CA:1175:G:H2'	1:CA:1176:A:C8	2.41	0.55
1:CA:1349:A:H2'	1:CA:1350:A:H8	1.72	0.55
1:CA:391:G:N1	1:CA:392:G:C5	2.74	0.55
2:CB:87:ARG:CZ	2:CB:233:SER:HB3	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:34:LEU:HD23	3:CC:34:LEU:C	2.27	0.55
3:CC:84:ILE:HD11	3:CC:88:ARG:NH2	2.20	0.55
4:CD:128:VAL:O	4:CD:130:GLY:N	2.38	0.55
5:CE:12:LEU:CD1	5:CE:31:LEU:HB3	2.34	0.55
8:CH:119:LEU:HD12	8:CH:124:ALA:N	2.20	0.55
9:CI:95:LYS:NZ	9:CI:96:LEU:HD13	2.20	0.55
11:CK:23:ALA:HA	11:CK:28:THR:HG23	1.87	0.55
12:CL:60:LEU:HD22	12:CL:60:LEU:H	1.71	0.55
12:CL:89:ARG:HB2	12:CL:89:ARG:HH11	1.70	0.55
14:CN:41:ARG:O	14:CN:44:LEU:HB3	2.06	0.55
15:CO:82:ILE:O	15:CO:82:ILE:HD13	2.06	0.55
16:CP:48:TRP:HE3	16:CP:49:LEU:H	1.54	0.55
16:CP:64:ALA:O	16:CP:65:GLN:C	2.45	0.55
35:DA:1326:U:O2'	35:DA:1327:C:H5'	2.06	0.55
35:DA:1361:G:O2'	35:DA:1362:C:H5'	2.07	0.55
35:DA:137:C:C2'	35:DA:139:G:H5'	2.36	0.55
35:DA:1612:C:C2'	35:DA:1613:G:O5'	2.55	0.55
35:DA:1652:A:N6	35:DA:1653:G:N2	2.54	0.55
35:DA:826:U:OP1	35:DA:2428:G:OP1	2.25	0.55
35:DA:270:A:O2'	35:DA:271:A:H5'	2.06	0.55
35:DA:2773:C:H2'	35:DA:2774:C:H6	1.70	0.55
35:DA:514:A:H2'	35:DA:515:A:C8	2.40	0.55
35:DA:601:C:O2	35:DA:605:C:H4'	2.05	0.55
35:DA:878:A:H3'	35:DA:879:G:C8	2.41	0.55
35:DA:995:C:C2	51:DU:57:PHE:CE2	2.94	0.55
39:DE:116:VAL:CG2	39:DE:117:MET:N	2.69	0.55
39:DE:57:LYS:HD3	39:DE:59:VAL:HG12	1.88	0.55
40:DF:110:LEU:O	40:DF:113:ALA:HB3	2.06	0.55
40:DF:41:LEU:O	40:DF:44:ARG:HG3	2.07	0.55
41:DG:40:ASN:H	41:DG:157:ILE:HA	1.71	0.55
44:DN:58:ASP:OD1	44:DN:124:ALA:HB1	2.05	0.55
44:DN:26:LEU:CD2	44:DN:30:ILE:HD11	2.37	0.55
35:DA:1245:G:C3'	46:DP:16:ARG:HH22	2.15	0.55
35:DA:662:G:O3'	46:DP:20:GLY:HA2	2.06	0.55
47:DQ:53:ALA:O	47:DQ:56:ARG:HB3	2.07	0.55
47:DQ:76:LYS:H	47:DQ:88:GLY:HA3	1.70	0.55
48:DR:55:ALA:HB1	48:DR:84:ALA:HB2	1.89	0.55
50:DT:65:LYS:HA	50:DT:65:LYS:HZ1	1.69	0.55
53:DW:80:PRO:HD2	53:DW:100:THR:HG21	1.89	0.55
55:DY:45:VAL:HG22	55:DY:62:GLU:CB	2.32	0.55
1:AA:1010:G:N2	1:AA:1020:U:H1'	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1285:A:H1'	1:AA:1286:A:OP2	2.07	0.55
1:AA:129(A):G:N2	1:AA:189(E):U:H1'	2.21	0.55
1:AA:709:G:H2'	1:AA:710:G:H8	1.72	0.55
1:AA:995:C:O2'	1:AA:996:A:H5'	2.06	0.55
2:AB:102:LEU:CD1	2:AB:102:LEU:N	2.69	0.55
2:AB:136:VAL:O	2:AB:140:HIS:HB2	2.07	0.55
2:AB:19:HIS:CD2	2:AB:20:GLU:HG2	2.42	0.55
2:AB:87:ARG:CZ	2:AB:233:SER:HB3	2.36	0.55
5:AE:101:ILE:CD1	5:AE:118:ILE:O	2.55	0.55
1:AA:1251:A:H5''	9:AI:12:GLU:OE1	2.06	0.55
13:AM:58:GLU:O	13:AM:62:ASN:HB3	2.06	0.55
15:AO:66:LEU:O	15:AO:69:TYR:HB3	2.07	0.55
27:B1:62:VAL:CG2	27:B1:67:ILE:HA	2.33	0.55
27:B1:74:VAL:O	27:B1:76:ARG:N	2.38	0.55
35:BA:1029:A:H2'	35:BA:1030:G:O4'	2.06	0.55
35:BA:1326:U:O2'	35:BA:1327:C:H5'	2.06	0.55
35:BA:141:A:H8	35:BA:1408:C:O2'	1.89	0.55
35:BA:1428:C:C4	35:BA:1569:A:H5''	2.41	0.55
35:BA:151:C:H42	35:BA:175:G:H1	1.52	0.55
35:BA:1792:G:O2'	35:BA:1793:C:H5'	2.07	0.55
35:BA:1902:C:H5'	38:BD:246:PRO:HD3	1.88	0.55
35:BA:2016:U:H2'	35:BA:2017:U:C6	2.41	0.55
35:BA:2465:C:O2'	35:BA:2466:C:H5'	2.05	0.55
35:BA:2570:G:H2'	35:BA:2571:C:H6	1.71	0.55
35:BA:300:A:H5''	55:BY:97:ARG:HH12	1.71	0.55
35:BA:401:A:H2'	35:BA:402:A:C8	2.41	0.55
35:BA:792:G:C4'	35:BA:793:A:H5'	2.36	0.55
35:BA:833:U:H5''	46:BP:48:PRO:HB2	1.87	0.55
36:BB:30:C:H2'	36:BB:31:C:O4'	2.07	0.55
41:BG:172:LEU:CA	41:BG:175:LEU:HD12	2.36	0.55
44:BN:26:LEU:HD11	44:BN:30:ILE:HD11	1.87	0.55
48:BR:28:LEU:HD13	48:BR:28:LEU:C	2.27	0.55
35:BA:1453:U:H5'	48:BR:63:ARG:HE	1.71	0.55
50:BT:83:ILE:HD11	50:BT:84:GLN:HE21	1.70	0.55
35:BA:1252:G:N3	51:BU:33:ARG:HD2	2.22	0.55
51:BU:79:PHE:O	51:BU:83:LEU:HD13	2.06	0.55
55:BY:28:LYS:HA	55:BY:39:VAL:N	2.20	0.55
56:BZ:36:LYS:O	56:BZ:37:VAL:HG13	2.06	0.55
1:CA:1321:C:H5''	1:CA:1322:C:H2'	1.88	0.55
1:CA:364:A:H2'	1:CA:365:U:O2	2.06	0.55
1:CA:783:C:H42	1:CA:800:G:N2	2.04	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:104:VAL:O	4:CD:107:ARG:HB2	2.07	0.55
4:CD:22:LYS:HB3	4:CD:22:LYS:NZ	2.20	0.55
12:CL:119:LYS:HD3	12:CL:120:TYR:HE1	1.71	0.55
13:CM:64:TRP:HE1	13:CM:66:LEU:HD12	1.72	0.55
3:CC:20:SER:O	14:CN:54:PRO:HG3	2.06	0.55
10:CJ:62:HIS:O	14:CN:59:ALA:HB3	2.06	0.55
16:CP:6:LEU:HD12	16:CP:6:LEU:N	2.20	0.55
17:CQ:9:VAL:HG12	17:CQ:10:VAL:N	2.21	0.55
18:CR:36:ASN:HB2	18:CR:39:VAL:HG21	1.88	0.55
23:CW:11:A:H2'	23:CW:12:G:H8	1.71	0.55
31:D5:40:LYS:HZ3	31:D5:50:GLY:HA2	1.70	0.55
35:DA:1029:A:H2'	35:DA:1030:G:O4'	2.07	0.55
35:DA:1175:U:H4'	35:DA:1176:G:C3'	2.37	0.55
35:DA:1297:C:H2'	35:DA:1298:C:C6	2.41	0.55
35:DA:1297:C:H2'	35:DA:1298:C:H6	1.71	0.55
35:DA:2590:A:O3'	38:DD:239:ARG:HG3	2.06	0.55
35:DA:2639:A:C3'	35:DA:2640:G:C5'	2.84	0.55
38:DD:143:HIS:CE1	38:DD:192:THR:HG23	2.42	0.55
38:DD:145:VAL:HB	38:DD:155:LEU:CB	2.37	0.55
39:DE:56:PRO:O	39:DE:57:LYS:O	2.25	0.55
40:DF:125:LEU:HB3	40:DF:196:LEU:CD2	2.37	0.55
40:DF:24:LEU:HB3	40:DF:25:PRO:HD2	1.88	0.55
40:DF:24:LEU:HB3	40:DF:25:PRO:CD	2.37	0.55
40:DF:32:LEU:HD23	40:DF:33:LEU:N	2.22	0.55
41:DG:99:MET:O	41:DG:103:LEU:HG	2.06	0.55
41:DG:120:LEU:HG	41:DG:179:PRO:HD2	1.88	0.55
41:DG:120:LEU:N	41:DG:179:PRO:O	2.40	0.55
44:DN:126:PRO:O	44:DN:127:ASP:OD1	2.24	0.55
45:DO:4:PRO:O	45:DO:5:GLN:HB3	2.05	0.55
47:DQ:109:VAL:HG13	47:DQ:113:GLN:OE1	2.07	0.55
47:DQ:114:ALA:O	47:DQ:116:GLU:N	2.39	0.55
48:DR:28:LEU:O	48:DR:30:THR:N	2.40	0.55
48:DR:73:VAL:HG23	48:DR:74:LYS:CD	2.35	0.55
48:DR:73:VAL:O	48:DR:76:VAL:HG12	2.07	0.55
51:DU:24:TYR:HB2	51:DU:29:SER:HB3	1.88	0.55
52:DV:2:PHE:CB	52:DV:42:GLY:HA2	2.30	0.55
56:DZ:11:GLU:H	56:DZ:11:GLU:CD	2.08	0.55
56:DZ:150:LEU:HD13	56:DZ:150:LEU:N	2.21	0.55
56:DZ:115:GLY:HA3	56:DZ:176:PRO:HA	1.88	0.55
1:AA:128:G:O2'	1:AA:129:U:H5'	2.07	0.55
1:AA:1360:A:H2'	1:AA:1361:G:O4'	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:17:U:H2'	1:AA:18:C:H6	1.68	0.55
1:AA:608:A:H4'	16:AP:32:TYR:OH	2.06	0.55
1:AA:783:C:H42	1:AA:800:G:N2	2.04	0.55
1:AA:971:G:H4'	1:AA:972:C:C5'	2.36	0.55
2:AB:99:GLY:O	2:AB:101:MET:N	2.39	0.55
3:AC:73:PRO:O	3:AC:75:VAL:N	2.39	0.55
5:AE:51:VAL:HB	5:AE:52:PRO:CD	2.35	0.55
7:AG:120:ILE:C	7:AG:124:LEU:HD12	2.27	0.55
9:AI:77:ILE:HG23	9:AI:81:ILE:HD11	1.89	0.55
14:AN:41:ARG:O	14:AN:44:LEU:HB3	2.07	0.55
20:AT:48:LYS:O	20:AT:49:ALA:HB2	2.06	0.55
25:AY:29:ARG:CZ	25:AY:110:ARG:HH21	2.19	0.55
27:B1:52:ARG:O	27:B1:53:VAL:HG12	2.06	0.55
31:B5:11:THR:HB	35:BA:1263:U:O3'	2.07	0.55
35:BA:1689:A:N6	35:BA:1698:A:H2	2.04	0.55
35:BA:1754:C:H2'	35:BA:1755:A:O4'	2.06	0.55
35:BA:212:G:O2'	35:BA:213:A:H5'	2.06	0.55
35:BA:499:U:O2'	35:BA:500:G:H5'	2.07	0.55
35:BA:926:A:H5'	35:BA:926:A:H8	1.70	0.55
35:BA:956:G:H22	35:BA:959:A:H3'	1.71	0.55
36:BB:40:U:N3	36:BB:43:C:H5''	2.22	0.55
30:B4:1:MET:H3	36:BB:43:C:H5'	1.71	0.55
37:BC:73:ARG:HG2	37:BC:92:ASP:OD2	2.07	0.55
37:BC:99:ILE:HG22	37:BC:99:ILE:O	2.06	0.55
38:BD:248:SER:HB2	38:BD:249:PRO:HD2	1.89	0.55
38:BD:34:VAL:CG2	38:BD:35:LYS:HZ2	2.19	0.55
39:BE:79:ARG:HH11	39:BE:79:ARG:HG2	1.72	0.55
40:BF:124:LEU:HD12	40:BF:125:LEU:N	2.22	0.55
40:BF:45:ARG:CG	40:BF:46:ARG:H	2.19	0.55
45:BO:86:ILE:CD1	45:BO:86:ILE:N	2.62	0.55
48:BR:9:LYS:NZ	48:BR:39:PRO:HA	2.21	0.55
48:BR:60:LEU:O	48:BR:61:HIS:C	2.43	0.55
35:BA:2377:A:H4'	49:BS:108:GLY:HA3	1.89	0.55
1:CA:1399:C:C2	1:CA:1401:G:C5	2.94	0.55
1:CA:1511:G:N1	1:CA:1525:G:C6	2.75	0.55
1:CA:472:A:C4'	16:CP:82:GLN:HE22	2.19	0.55
1:CA:540:G:C2	1:CA:541:G:C4	2.95	0.55
2:CB:124:SER:OG	2:CB:125:PRO:HD2	2.07	0.55
2:CB:82:ARG:HG2	2:CB:82:ARG:HH11	1.71	0.55
4:CD:62:GLN:HB3	4:CD:66:ARG:CZ	2.37	0.55
7:CG:135:VAL:HG12	7:CG:139:GLU:HG3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:50:ILE:O	7:CG:54:THR:O	2.24	0.55
7:CG:75:VAL:O	7:CG:75:VAL:HG23	2.06	0.55
9:CI:4:TYR:CB	9:CI:19:LEU:HB2	2.35	0.55
11:CK:43:SER:OG	11:CK:47:VAL:HG11	2.07	0.55
12:CL:70:ILE:N	12:CL:70:ILE:CD1	2.68	0.55
18:CR:25:THR:O	18:CR:26:LEU:HG	2.07	0.55
27:D1:56:GLN:O	27:D1:57:GLU:HB2	2.06	0.55
27:D1:85:LEU:O	27:D1:86:SER:C	2.45	0.55
35:DA:1196:C:H2'	35:DA:1197:G:C8	2.41	0.55
35:DA:1446:C:N4	35:DA:1465:G:H1	2.02	0.55
35:DA:2537:U:H2'	35:DA:2538:C:C6	2.42	0.55
35:DA:2559:C:O2	35:DA:2559:C:H2'	2.06	0.55
35:DA:2643:G:C2	35:DA:2772:C:N3	2.74	0.55
35:DA:2809:A:O2'	35:DA:2810:A:H5'	2.06	0.55
35:DA:302:C:O2'	35:DA:303:U:H5'	2.07	0.55
36:DB:7:G:H4'	49:DS:29:PHE:HD2	1.65	0.55
37:DC:41:VAL:HB	37:DC:178:ALA:HB1	1.88	0.55
38:DD:143:HIS:HB3	38:DD:194:GLY:O	2.05	0.55
38:DD:218:ARG:HB3	38:DD:219:PRO:HD2	1.88	0.55
39:DE:2:LYS:CD	39:DE:95:ILE:HG22	2.36	0.55
39:DE:88:GLY:O	39:DE:89:ASP:HB2	2.07	0.55
40:DF:139:PHE:CB	40:DF:166:ALA:HB1	2.36	0.55
41:DG:137:GLU:CG	41:DG:138:GLN:N	2.64	0.55
42:DH:19:VAL:HG11	42:DH:44:VAL:HG22	1.88	0.55
45:DO:16:ALA:CB	45:DO:43:VAL:HG13	2.34	0.55
46:DP:95:VAL:O	46:DP:125:VAL:HG23	2.07	0.55
51:DU:92:ARG:C	51:DU:94:ASN:N	2.58	0.55
52:DV:27:ALA:O	52:DV:29:PRO:N	2.39	0.55
51:DU:90:VAL:CG2	52:DV:39:LEU:HD12	2.36	0.55
28:D2:26:ARG:CZ	54:DX:5:TYR:HB3	2.36	0.55
1:AA:1031:G:H2'	1:AA:1032:G:C8	2.41	0.55
1:AA:1288:A:O2'	1:AA:1289:A:H5'	2.07	0.55
1:AA:450:G:H4'	16:AP:41:PRO:O	2.07	0.55
1:AA:453:A:H2'	1:AA:454:C:C6	2.42	0.55
1:AA:781:A:H2'	1:AA:782:A:H5'	1.88	0.55
1:AA:788:U:C5	1:AA:789:U:H5	2.24	0.55
1:AA:857:C:H2'	1:AA:858:G:O4'	2.05	0.55
7:AG:75:VAL:HG23	7:AG:75:VAL:O	2.06	0.55
11:AK:83:ILE:HA	11:AK:109:VAL:O	2.06	0.55
11:AK:57:THR:HG23	11:AK:58:PRO:HD2	1.89	0.55
12:AL:102:ARG:CG	12:AL:102:ARG:HH11	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:10:VAL:HG12	17:AQ:53:LEU:HD12	1.87	0.55
18:AR:35:ARG:C	18:AR:37:VAL:N	2.59	0.55
25:AY:144:ALA:HB2	25:AY:149:LEU:HD12	1.89	0.55
25:AY:37:LEU:C	25:AY:37:LEU:HD12	2.27	0.55
31:B5:26:THR:O	31:B5:26:THR:HG23	2.06	0.55
32:B6:10:LEU:HD22	32:B6:10:LEU:N	2.22	0.55
33:B7:30:VAL:O	33:B7:31:LEU:C	2.44	0.55
35:BA:1515:G:H2'	35:BA:1516:C:H5'	1.88	0.55
35:BA:1649:G:O2'	35:BA:1650:G:H5'	2.06	0.55
35:BA:2720:U:O2	35:BA:2720:U:C2'	2.54	0.55
35:BA:45:C:H2'	35:BA:47:C:C6	2.42	0.55
38:BD:257:LEU:CD2	38:BD:257:LEU:C	2.74	0.55
35:BA:2632:A:H1'	39:BE:61:ARG:NH1	2.22	0.55
35:BA:607:U:OP1	40:BF:103:LYS:N	2.39	0.55
44:BN:43:THR:O	44:BN:44:PRO:C	2.43	0.55
45:BO:2:ILE:CD1	45:BO:6:THR:HG21	2.36	0.55
46:BP:100:LEU:CD2	46:BP:100:LEU:H	2.20	0.55
35:BA:662:G:O3'	46:BP:20:GLY:HA2	2.06	0.55
47:BQ:114:ALA:O	47:BQ:116:GLU:N	2.40	0.55
53:BW:40:ASN:O	53:BW:41:LYS:HG2	2.07	0.55
56:BZ:125:LEU:O	56:BZ:126:VAL:HG22	2.07	0.55
56:BZ:114:GLY:CA	56:BZ:177:PRO:HB3	2.31	0.55
56:BZ:5:LEU:HD11	56:BZ:43:GLU:O	2.05	0.55
1:CA:1230:C:O2'	1:CA:1231:G:H5'	2.07	0.55
1:CA:1256:A:H61	1:CA:1278:U:H1'	1.71	0.55
1:CA:1296:C:H3'	1:CA:1297:C:H6	1.72	0.55
1:CA:383:A:H2'	1:CA:384:G:C5'	2.28	0.55
1:CA:452:A:C2	1:CA:453:A:C4	2.94	0.55
1:CA:656:C:H4'	15:CO:62:GLN:HE22	1.71	0.55
2:CB:105:PHE:HA	2:CB:108:ILE:HG22	1.87	0.55
6:CF:69:GLU:H	6:CF:69:GLU:CD	2.10	0.55
11:CK:30:VAL:HG21	11:CK:65:ALA:HA	1.89	0.55
13:CM:58:GLU:O	13:CM:62:ASN:HB3	2.05	0.55
13:CM:70:LEU:O	13:CM:74:VAL:HG23	2.07	0.55
13:CM:76:ALA:HA	13:CM:79:LYS:HD2	1.89	0.55
25:CY:162:GLN:O	25:CY:163:LYS:C	2.45	0.55
31:D5:20:ARG:HA	31:D5:23:HIS:HD2	1.70	0.55
35:DA:1331:A:H2'	35:DA:1333:C:H5	1.71	0.55
35:DA:1415:U:O2	35:DA:1415:U:H2'	2.06	0.55
35:DA:1899:G:H22	35:DA:1902:C:N4	2.04	0.55
35:DA:2079:U:H3	35:DA:2241:A:N6	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2308:G:O6	35:DA:2310:A:H2'	2.07	0.55
35:DA:2460:U:O2'	35:DA:2461:C:H5'	2.06	0.55
35:DA:268:C:C2'	35:DA:268:C:O2	2.53	0.55
35:DA:470:A:OP1	40:DF:59:TYR:HE2	1.89	0.55
35:DA:549:G:H2'	35:DA:551:G:C4'	2.36	0.55
35:DA:66:C:C2'	35:DA:67:U:H5'	2.37	0.55
35:DA:913:U:H4'	35:DA:914:C:OP1	2.06	0.55
38:DD:260:ARG:NH2	38:DD:264:LYS:HD3	2.21	0.55
40:DF:74:ARG:O	40:DF:75:HIS:ND1	2.40	0.55
40:DF:89:VAL:CG1	40:DF:90:PHE:H	2.20	0.55
41:DG:17:PRO:C	41:DG:19:LEU:H	2.10	0.55
41:DG:38:VAL:HA	41:DG:92:VAL:O	2.07	0.55
42:DH:97:ARG:HD2	42:DH:104:GLU:OE1	2.07	0.55
42:DH:65:HIS:ND1	42:DH:66:GLY:N	2.55	0.55
44:DN:46:VAL:HG21	44:DN:48:MET:HG3	1.88	0.55
46:DP:46:LYS:CB	46:DP:52:GLU:HG2	2.37	0.55
48:DR:2:ARG:HD2	48:DR:2:ARG:C	2.26	0.55
56:DZ:17:ALA:O	56:DZ:18:LEU:C	2.45	0.55
1:AA:115:G:O2'	1:AA:116:A:OP2	2.22	0.55
1:AA:1439:C:OP1	20:AT:38:LYS:HD2	2.06	0.55
1:AA:231:G:O2'	1:AA:232:G:H5'	2.06	0.55
1:AA:357:G:OP1	1:AA:366:C:O2'	2.24	0.55
1:AA:475:G:H2'	1:AA:476:G:H8	1.71	0.55
1:AA:792:A:H4'	1:AA:793:U:O5'	2.07	0.55
1:AA:856:C:H2'	1:AA:857:C:H6	1.72	0.55
1:AA:980:C:H5'	1:AA:981:U:C5	2.41	0.55
3:AC:34:LEU:HD23	3:AC:34:LEU:C	2.27	0.55
12:AL:89:ARG:O	12:AL:89:ARG:HD3	2.07	0.55
13:AM:91:ARG:HG3	13:AM:98:VAL:HG13	1.89	0.55
14:AN:51:GLY:C	14:AN:53:LEU:H	2.10	0.55
19:AS:6:LYS:CD	19:AS:6:LYS:H	2.20	0.55
20:AT:11:SER:HA	20:AT:13:LEU:HD12	1.88	0.55
22:AV:28:G:H2'	22:AV:29:G:H8	1.72	0.55
35:BA:1114:G:H2'	35:BA:1115:G:C5'	2.32	0.55
35:BA:1131:G:N3	35:BA:1132:A:C8	2.74	0.55
35:BA:585:G:H2'	35:BA:1251:C:N4	2.22	0.55
35:BA:1352:U:O2'	35:BA:1353:A:H5'	2.07	0.55
35:BA:1496:A:H2'	35:BA:1498:C:C5	2.42	0.55
35:BA:1786:A:N1	35:BA:2606:C:H1'	2.22	0.55
35:BA:201:C:C2'	35:BA:202:U:H5'	2.37	0.55
35:BA:247:G:H4'	35:BA:386:G:C5	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2547:U:O2'	35:BA:2548:G:H5'	2.07	0.55
35:BA:2801:A:H4'	35:BA:2801(A):A:O5'	2.07	0.55
35:BA:2801:A:O2'	35:BA:2895:U:H4'	2.06	0.55
35:BA:2852:G:H2'	35:BA:2853:C:H6	1.71	0.55
35:BA:778:G:H5''	38:BD:48:ARG:HD2	1.88	0.55
38:BD:146:GLU:CA	38:BD:153:ALA:HA	2.33	0.55
38:BD:259:THR:O	38:BD:260:ARG:C	2.45	0.55
39:BE:173:VAL:HG12	39:BE:174:ASP:N	2.20	0.55
41:BG:141:PHE:O	41:BG:144:ILE:HG22	2.07	0.55
41:BG:96:ARG:HA	41:BG:99:MET:CE	2.36	0.55
34:B8:59:LYS:HG3	46:BP:49:ARG:HD2	1.89	0.55
46:BP:56:SER:C	46:BP:58:THR:H	2.10	0.55
53:BW:64:MET:HE2	53:BW:109:GLU:HG3	1.88	0.55
1:CA:1202:G:C2'	1:CA:1203:C:H5'	2.36	0.55
1:CA:1325:C:H2'	1:CA:1326:C:H5'	1.88	0.55
1:CA:1463:C:H2'	1:CA:1464:G:O4'	2.06	0.55
1:CA:437:U:O2'	1:CA:438:G:H5'	2.06	0.55
1:CA:44:G:C2	1:CA:45:U:H1'	2.42	0.55
5:CE:72:GLN:O	5:CE:73:ASN:HB3	2.05	0.55
9:CI:117:HIS:HB2	9:CI:121:ARG:HD2	1.89	0.55
10:CJ:32:ALA:HB1	10:CJ:75:ILE:HG13	1.89	0.55
11:CK:65:ALA:HB3	11:CK:97:ALA:CB	2.36	0.55
11:CK:80:VAL:HG23	11:CK:80:VAL:O	2.06	0.55
11:CK:96:ARG:CA	11:CK:99:GLN:HG2	2.36	0.55
6:CF:100:ASN:H	18:CR:23:LYS:NZ	2.05	0.55
19:CS:22:LEU:HD22	19:CS:27:GLU:H	1.70	0.55
20:CT:30:LYS:HZ3	20:CT:72:LEU:HD21	1.72	0.55
26:D0:23:VAL:HG12	26:D0:24:LYS:N	2.22	0.55
35:DA:121:G:H2'	35:DA:122:G:H8	1.71	0.55
35:DA:1445(A):C:H2'	35:DA:1446:C:H6	1.70	0.55
35:DA:1931:U:O2'	35:DA:1932:A:H5'	2.06	0.55
35:DA:2098:U:H2'	35:DA:2099:U:C6	2.42	0.55
35:DA:2720:U:O2	35:DA:2720:U:C2'	2.54	0.55
35:DA:2732:G:H3'	35:DA:2733:A:C5'	2.36	0.55
35:DA:2801(A):A:H5'	35:DA:2802:G:C8	2.41	0.55
35:DA:2888:C:H2'	35:DA:2889:C:H6	1.70	0.55
35:DA:468:G:H2'	35:DA:469:G:O4'	2.06	0.55
35:DA:79:G:O2'	35:DA:80:G:H5'	2.06	0.55
35:DA:921:G:H2'	35:DA:922:U:C6	2.42	0.55
39:DE:2:LYS:CE	39:DE:95:ILE:HG22	2.37	0.55
39:DE:64:LYS:C	39:DE:66:HIS:N	2.59	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DF:124:LEU:HD12	40:DF:125:LEU:N	2.21	0.55
40:DF:170:LEU:HD21	40:DF:172:TRP:CE2	2.42	0.55
40:DF:45:ARG:CG	40:DF:46:ARG:H	2.19	0.55
43:DI:4:ILE:C	43:DI:5:LEU:HD23	2.27	0.55
43:DI:94:ALA:C	43:DI:96:ASP:N	2.59	0.55
44:DN:43:THR:O	44:DN:45:ASN:N	2.40	0.55
44:DN:75:TYR:O	44:DN:82:LEU:O	2.25	0.55
47:DQ:38:GLU:HB2	47:DQ:127:ILE:CG2	2.37	0.55
48:DR:74:LYS:HA	48:DR:77:ARG:HD2	1.88	0.55
35:DA:2377:A:H4'	49:DS:108:GLY:HA3	1.88	0.55
49:DS:27:SER:N	49:DS:38:GLN:O	2.39	0.55
49:DS:40:ILE:HG23	49:DS:46:VAL:O	2.07	0.55
49:DS:99:LYS:O	49:DS:101:LEU:N	2.40	0.55
50:DT:13:ARG:HH12	50:DT:15:VAL:HG12	1.68	0.55
50:DT:62:THR:CG2	50:DT:75:ILE:HG13	2.37	0.55
44:DN:40:PRO:CB	51:DU:64:ARG:HH22	2.19	0.55
56:DZ:10:ARG:HG3	56:DZ:38:TYR:H	1.72	0.55
1:AA:1416:G:O2'	1:AA:1417:G:H5'	2.07	0.55
1:AA:363:A:C5	12:AL:31:PRO:HD2	2.40	0.55
1:AA:666:G:H1'	1:AA:741:G:N2	2.21	0.55
1:AA:853:G:H2'	1:AA:854:G:H8	1.71	0.55
2:AB:36:ARG:NE	2:AB:37:ASN:H	2.05	0.55
2:AB:61:LEU:O	2:AB:64:ARG:HG2	2.05	0.55
3:AC:188:LEU:HD22	3:AC:188:LEU:N	2.21	0.55
4:AD:156:GLU:O	4:AD:159:ARG:HB2	2.07	0.55
4:AD:93:PHE:CE1	4:AD:97:LEU:HD12	2.42	0.55
1:AA:972:C:H2'	10:AJ:55:LYS:HD3	1.88	0.55
12:AL:9:GLN:O	12:AL:10:LEU:C	2.45	0.55
13:AM:93:ARG:HA	13:AM:93:ARG:HE	1.72	0.55
16:AP:82:GLN:H	16:AP:82:GLN:HE21	1.38	0.55
17:AQ:80:GLY:O	17:AQ:81:ARG:HG2	2.06	0.55
22:AV:40:C:O2'	22:AV:41:C:H5'	2.07	0.55
27:B1:83:GLU:O	27:B1:85:LEU:N	2.30	0.55
28:B2:41:ILE:O	28:B2:43:GLN:N	2.40	0.55
31:B5:29:THR:O	31:B5:42:PRO:HD3	2.07	0.55
31:B5:46:CYS:SG	31:B5:48:GLU:HG3	2.47	0.55
35:BA:81:G:H1	35:BA:105:C:H42	1.55	0.55
35:BA:1335:U:H2'	35:BA:1336:A:H8	1.72	0.55
35:BA:2476:A:N3	35:BA:2477:C:H5'	2.22	0.55
35:BA:2661:G:H2'	35:BA:2662:A:H8	1.69	0.55
35:BA:2863:C:C3'	35:BA:2864:G:H5''	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:493:G:H3'	35:BA:494:G:H5''	1.89	0.55
35:BA:565:C:H2'	35:BA:566:U:H6	1.72	0.55
35:BA:662:G:O2'	35:BA:663:G:H5'	2.07	0.55
35:BA:893:C:H5	35:BA:894:C:C2	2.25	0.55
35:BA:863:A:H4'	36:BB:101:G:N2	2.21	0.55
43:BI:10:GLU:CD	43:BI:11:ASN:H	2.10	0.55
44:BN:128:HIS:O	44:BN:130:HIS:N	2.40	0.55
44:BN:46:VAL:O	44:BN:47:ALA:HB2	2.06	0.55
45:BO:87:ILE:HG23	45:BO:88:ASN:N	2.22	0.55
49:BS:99:LYS:O	49:BS:101:LEU:N	2.40	0.55
50:BT:28:VAL:HB	50:BT:88:ILE:HG12	1.89	0.55
1:CA:1031:G:H2'	1:CA:1032:G:C8	2.42	0.55
1:CA:1067:A:H8	1:CA:1067:A:O5'	1.89	0.55
1:CA:11:G:C5	1:CA:12:U:C5	2.95	0.55
1:CA:1285:A:H1'	1:CA:1286:A:OP2	2.07	0.55
1:CA:128:G:O2'	1:CA:129:U:H5'	2.06	0.55
1:CA:1424:C:C2	1:CA:1425:U:C6	2.95	0.55
1:CA:197:A:N3	1:CA:198:G:H1'	2.22	0.55
1:CA:498:U:C2'	1:CA:498:U:O2	2.55	0.55
1:CA:555:C:H2'	1:CA:556:C:C6	2.42	0.55
1:CA:817:C:N4	1:CA:1529:G:H1	2.05	0.55
3:CC:188:LEU:HD22	3:CC:188:LEU:N	2.22	0.55
6:CF:68:PRO:CG	6:CF:71:ARG:HE	2.18	0.55
9:CI:47:LEU:N	9:CI:47:LEU:HD12	2.22	0.55
11:CK:58:PRO:HD3	11:CK:89:ALA:CB	2.35	0.55
19:CS:15:LEU:O	19:CS:19:VAL:HG23	2.07	0.55
25:CY:15:GLN:CA	25:CY:168:PHE:HZ	2.19	0.55
33:D7:30:VAL:O	33:D7:31:LEU:C	2.43	0.55
35:DA:1572:A:O2'	35:DA:1573:G:H5'	2.07	0.55
35:DA:1830:C:N4	35:DA:1975:G:H1	2.02	0.55
35:DA:1999:C:H2'	35:DA:2000:G:C8	2.39	0.55
35:DA:1998:G:O2'	35:DA:1999:C:H5'	2.07	0.55
35:DA:2345:G:H5''	35:DA:2347:C:O4'	2.07	0.55
35:DA:2785:C:H2'	35:DA:2786:U:C6	2.41	0.55
35:DA:285:C:H2'	35:DA:286:C:C5'	2.36	0.55
27:D1:69:LYS:NZ	35:DA:372:G:P	2.80	0.55
35:DA:8:A:H5''	44:DN:51:PHE:HZ	1.72	0.55
35:DA:848:G:N3	35:DA:933:A:H1'	2.21	0.55
38:DD:160:GLY:N	38:DD:196:VAL:HB	2.22	0.55
39:DE:133:LYS:N	39:DE:134:ILE:HD13	2.22	0.55
41:DG:102:PHE:HA	41:DG:105:LYS:HZ1	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:9:ARG:O	41:DG:13:GLU:HG2	2.07	0.55
41:DG:145:THR:OG1	41:DG:148:MET:HB3	2.07	0.55
41:DG:166:ASP:OD1	41:DG:170:ARG:HD2	2.07	0.55
41:DG:170:ARG:HG3	41:DG:170:ARG:HH11	1.71	0.55
41:DG:22:ARG:HH11	41:DG:22:ARG:HG2	1.72	0.55
43:DI:83:ALA:HA	43:DI:89:TYR:HD1	1.72	0.55
43:DI:87:LYS:CE	43:DI:121:LYS:HG2	2.36	0.55
54:DX:33:LYS:C	54:DX:35:THR:H	2.09	0.55
55:DY:16:ALA:CA	55:DY:21:LYS:HD2	2.36	0.55
55:DY:31:LEU:HB2	55:DY:36:ALA:O	2.07	0.55
56:DZ:40:ASP:O	56:DZ:44:PHE:HB2	2.05	0.55
2:AB:80:ILE:HD11	2:AB:215:LEU:HD12	1.88	0.55
5:AE:131:ILE:H	5:AE:131:ILE:CD1	2.16	0.55
8:AH:45:ILE:HB	8:AH:62:TYR:O	2.06	0.55
10:AJ:16:LEU:HD22	10:AJ:19:SER:OG	2.06	0.55
13:AM:92:HIS:CE1	13:AM:98:VAL:HG21	2.41	0.55
14:AN:53:LEU:HB3	14:AN:56:VAL:HG21	1.89	0.55
16:AP:8:ARG:NH2	16:AP:15:PRO:HG3	2.22	0.55
17:AQ:31:LEU:HG	17:AQ:32:TYR:CE2	2.42	0.55
17:AQ:67:LYS:CA	17:AQ:70:ARG:HH12	2.19	0.55
18:AR:75:ILE:HG22	18:AR:75:ILE:O	2.06	0.55
25:AY:14:MET:CE	25:AY:165:THR:HG23	2.37	0.55
25:AY:78:ALA:HA	25:AY:81:LYS:HD2	1.89	0.55
27:B1:87:PRO:N	27:B1:89:GLU:HG2	2.20	0.55
27:B1:87:PRO:HB2	27:B1:91:LYS:CE	2.37	0.55
34:B8:48:PHE:N	34:B8:48:PHE:CD1	2.73	0.55
34:B8:55:ALA:O	34:B8:59:LYS:NZ	2.36	0.55
35:BA:1277:G:O2'	35:BA:1278:A:H5'	2.06	0.55
35:BA:1417:C:C2'	35:BA:1418:G:H5'	2.37	0.55
35:BA:143(A):C:H2'	35:BA:143(A):C:O2	2.06	0.55
35:BA:1590:U:H2'	35:BA:1591:G:C5'	2.28	0.55
35:BA:2718:G:H2'	35:BA:2719:G:C8	2.42	0.55
35:BA:2781:A:H5'	35:BA:2782:G:H5'	1.87	0.55
35:BA:685:A:C2	35:BA:787:U:H1'	2.42	0.55
35:BA:2086:U:OP1	38:BD:262:ARG:HG2	2.07	0.55
39:BE:55:ASN:ND2	39:BE:75:VAL:HG13	2.22	0.55
40:BF:24:LEU:HB3	40:BF:25:PRO:HD2	1.88	0.55
40:BF:65:TRP:O	40:BF:67:GLN:N	2.39	0.55
41:BG:115:ARG:NH2	41:BG:136:ARG:HD2	2.19	0.55
41:BG:178:PHE:CB	41:BG:180:PHE:HE1	2.14	0.55
44:BN:78:TYR:CD1	44:BN:79:PRO:HD3	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BN:75:TYR:O	44:BN:82:LEU:O	2.24	0.55
44:BN:82:LEU:HD12	44:BN:83:LYS:N	2.21	0.55
44:BN:99:LEU:O	44:BN:103:VAL:HG23	2.06	0.55
50:BT:110:ILE:HG22	50:BT:111:ARG:N	2.22	0.55
51:BU:91:ASP:O	51:BU:92:ARG:O	2.25	0.55
51:BU:99:ALA:HB2	51:BU:106:PHE:CD1	2.42	0.55
52:BV:96:ILE:HG23	52:BV:97:LYS:N	2.22	0.55
35:BA:310:A:OP1	55:BY:17:SER:O	2.25	0.55
55:BY:86:ARG:HB3	55:BY:88:LYS:NZ	2.20	0.55
56:BZ:104:PHE:HA	56:BZ:139:VAL:HB	1.89	0.55
1:CA:788:U:C4	1:CA:789:U:C5	2.95	0.55
2:CB:16:HIS:HA	2:CB:210:SER:HB2	1.87	0.55
4:CD:133:VAL:CG1	4:CD:135:LEU:H	2.20	0.55
5:CE:15:ARG:HG2	5:CE:26:PHE:CD2	2.42	0.55
6:CF:27:GLN:NE2	6:CF:27:GLN:HA	2.21	0.55
6:CF:52:ILE:O	6:CF:86:ARG:NH1	2.40	0.55
9:CI:7:THR:HB	9:CI:83:ARG:HH11	1.72	0.55
15:CO:64:ARG:O	15:CO:65:ARG:C	2.45	0.55
16:CP:18:ARG:O	16:CP:20:VAL:HG12	2.07	0.55
25:CY:40:HIS:O	25:CY:41:LEU:C	2.45	0.55
33:D7:8:ASN:ND2	33:D7:8:ASN:C	2.60	0.55
34:D8:62:LEU:N	34:D8:63:PRO:HD2	2.21	0.55
35:DA:2513:G:H1	35:DA:2571:C:H42	1.54	0.55
35:DA:2677:G:C4	35:DA:2678:C:C5	2.95	0.55
35:DA:2679:A:O2'	35:DA:2680:C:H5'	2.07	0.55
35:DA:738:G:C6	35:DA:739:G:C2	2.95	0.55
35:DA:874:G:H2'	35:DA:875:G:H8	1.71	0.55
35:DA:952:G:C6	35:DA:953:A:N7	2.75	0.55
38:DD:35:LYS:HD3	38:DD:63:ARG:CB	2.23	0.55
39:DE:110:GLY:O	48:DR:2:ARG:CZ	2.55	0.55
40:DF:114:VAL:CG2	40:DF:115:ALA:H	2.14	0.55
41:DG:160:VAL:HG13	41:DG:161:THR:N	2.21	0.55
41:DG:61:ALA:HB1	41:DG:68:PRO:HD3	1.89	0.55
42:DH:153:LYS:N	42:DH:153:LYS:HD3	2.22	0.55
45:DO:104:ARG:O	45:DO:107:ARG:N	2.38	0.55
35:DA:598:G:H5'	46:DP:15:ARG:CB	2.37	0.55
50:DT:62:THR:HB	50:DT:74:ARG:O	2.07	0.55
52:DV:72:VAL:CA	52:DV:88:ARG:HH12	2.17	0.55
54:DX:70:LEU:HG	54:DX:71:GLY:N	2.20	0.55
55:DY:39:VAL:O	55:DY:40:GLU:HG2	2.07	0.55
56:DZ:143:GLY:O	56:DZ:144:LEU:HD22	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1018:C:H2'	1:AA:1019:C:C6	2.42	0.55
1:AA:11:G:C5	1:AA:12:U:C5	2.95	0.55
1:AA:872:A:C2	1:AA:874:G:C6	2.95	0.55
1:AA:965:A:C2	1:AA:969:A:C2	2.95	0.55
2:AB:54:THR:HG22	2:AB:58:ILE:CD1	2.37	0.55
3:AC:84:ILE:HD11	3:AC:88:ARG:NH2	2.22	0.55
4:AD:65:ARG:HB2	4:AD:75:PHE:CZ	2.42	0.55
5:AE:133:TYR:H	5:AE:133:TYR:HD1	1.53	0.55
11:AK:34:ASP:HB3	11:AK:40:ILE:HD11	1.88	0.55
11:AK:30:VAL:HG21	11:AK:65:ALA:HA	1.87	0.55
12:AL:100:ILE:O	12:AL:102:ARG:N	2.40	0.55
27:B1:18:ILE:O	27:B1:18:ILE:HG22	2.07	0.55
31:B5:22:HIS:N	31:B5:22:HIS:ND1	2.54	0.55
35:BA:1160:G:N2	52:BV:10:LYS:HE3	2.22	0.55
35:BA:1196:C:H2'	35:BA:1197:G:H8	1.69	0.55
35:BA:1275:A:C4	48:BR:16:HIS:CE1	2.95	0.55
35:BA:2202:C:C2'	38:BD:151:LYS:HZ1	2.20	0.55
23:AW:77:A:C2	35:BA:2421:G:C6	2.95	0.55
35:BA:285:C:H2'	35:BA:286:C:C5'	2.36	0.55
35:BA:464:U:H2'	35:BA:465:G:O4'	2.07	0.55
35:BA:627:A:H4'	35:BA:628:G:OP1	2.06	0.55
41:BG:36:LYS:HG2	41:BG:37:VAL:N	2.21	0.55
42:BH:157:TYR:HD1	42:BH:170:ARG:O	1.88	0.55
44:BN:126:PRO:O	44:BN:127:ASP:OD1	2.25	0.55
44:BN:41:ASP:N	51:BU:64:ARG:NH1	2.51	0.55
44:BN:46:VAL:HG21	44:BN:48:MET:HG3	1.89	0.55
44:BN:72:TYR:HB3	44:BN:74:ARG:HG2	1.89	0.55
35:BA:906:G:H5'	47:BQ:26:TYR:OH	2.07	0.55
39:BE:12:THR:HG23	50:BT:8:LYS:HE2	1.88	0.55
51:BU:7:GLY:O	51:BU:8:VAL:HG22	2.07	0.55
54:BX:12:VAL:HG11	54:BX:27:THR:CG2	2.35	0.55
56:BZ:42:VAL:HG13	56:BZ:43:GLU:H	1.70	0.55
56:BZ:56:VAL:CG1	56:BZ:57:ILE:N	2.69	0.55
56:BZ:61:LEU:O	56:BZ:63:ASP:N	2.35	0.55
1:CA:192:U:H2'	1:CA:193:C:C6	2.41	0.55
1:CA:533:A:H1'	1:CA:534:U:OP1	2.06	0.55
1:CA:769:G:H2'	1:CA:770:C:H6	1.72	0.55
2:CB:75:LYS:HE3	2:CB:75:LYS:HA	1.89	0.55
11:CK:83:ILE:HA	11:CK:109:VAL:O	2.06	0.55
16:CP:19:ILE:HG22	16:CP:36:ILE:CG1	2.36	0.55
19:CS:64:GLU:HG3	19:CS:65:ASN:N	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:D8:39:LYS:HG2	34:D8:42:ARG:HH11	1.71	0.55
34:D8:53:PRO:C	34:D8:55:ALA:H	2.10	0.55
35:DA:1021:A:C8	35:DA:1021:A:H3'	2.41	0.55
35:DA:1668:A:N7	35:DA:1674:G:C6	2.75	0.55
35:DA:2055:C:H4'	35:DA:2056:G:H5''	1.89	0.55
35:DA:215:G:H4'	35:DA:216:A:O5'	2.06	0.55
35:DA:2223:G:H2'	35:DA:2224:G:H5'	1.88	0.55
35:DA:271(G):C:O2'	35:DA:271(H):G:H5'	2.07	0.55
35:DA:782:A:C2	38:DD:226:MET:CG	2.81	0.55
35:DA:991:C:H42	35:DA:1163:G:H1	1.54	0.55
37:DC:73:ARG:HG2	37:DC:92:ASP:OD2	2.06	0.55
38:DD:65:ILE:O	38:DD:65:ILE:HD12	2.07	0.55
39:DE:36:ARG:NH2	39:DE:88:GLY:HA3	2.22	0.55
42:DH:96:ALA:HB2	42:DH:105:LEU:HB3	1.89	0.55
44:DN:43:THR:O	44:DN:44:PRO:C	2.45	0.55
47:DQ:32:TYR:HD1	47:DQ:32:TYR:H	1.54	0.55
48:DR:85:PRO:O	48:DR:87:TYR:N	2.40	0.55
50:DT:89:VAL:CG1	50:DT:91:ARG:HE	2.20	0.55
51:DU:95:LEU:HD12	52:DV:11:GLN:HG3	1.89	0.55
52:DV:61:VAL:HG23	52:DV:100:ARG:N	2.21	0.55
52:DV:14:VAL:HG11	52:DV:98:GLU:HG3	1.88	0.55
54:DX:52:VAL:O	54:DX:53:LYS:CB	2.55	0.55
55:DY:47:LYS:HG3	55:DY:60:PHE:CE2	2.41	0.55
1:AA:1053:G:C6	1:AA:1199:U:H2'	2.42	0.54
1:AA:1203:C:H2'	1:AA:1204:A:C8	2.41	0.54
1:AA:472:A:C4'	16:AP:82:GLN:HE22	2.19	0.54
1:AA:865:A:H5'	1:AA:1078:U:H5	1.72	0.54
1:AA:78:G:H22	1:AA:91:C:H42	1.53	0.54
2:AB:159:PRO:C	2:AB:161:ALA:N	2.57	0.54
2:AB:212:GLN:HG3	2:AB:235:SER:HB2	1.89	0.54
3:AC:86:VAL:O	3:AC:89:GLU:HB3	2.06	0.54
4:AD:62:GLN:HB3	4:AD:66:ARG:NH1	2.22	0.54
5:AE:80:ILE:HD11	5:AE:91:LEU:HB2	1.89	0.54
8:AH:129:VAL:HG23	8:AH:130:GLY:N	2.16	0.54
11:AK:85:ARG:HG2	11:AK:111:ASP:O	2.06	0.54
18:AR:56:THR:HG21	18:AR:63:GLN:HE22	1.72	0.54
35:BA:1246:A:OP1	46:BP:18:ARG:HG3	2.07	0.54
35:BA:1415:U:O2	35:BA:1415:U:H2'	2.06	0.54
35:BA:1790:C:H2'	35:BA:1791:A:C8	2.41	0.54
35:BA:201:C:O2'	35:BA:202:U:H5'	2.07	0.54
35:BA:2052:G:C2	39:BE:149:ARG:HA	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2248:C:H3'	35:BA:2249:U:H6	1.72	0.54
35:BA:2284:C:H42	35:BA:2384:G:H1	1.54	0.54
35:BA:2571:C:O2	35:BA:2571:C:H2'	2.06	0.54
35:BA:2513:G:H1	35:BA:2571:C:H42	1.55	0.54
35:BA:2022:U:O2'	35:BA:2617:C:H5'	2.07	0.54
35:BA:360:G:O2'	35:BA:361:G:H5'	2.07	0.54
35:BA:409:C:O2'	35:BA:410:G:H5'	2.06	0.54
35:BA:535:C:O2'	35:BA:536:A:H5'	2.07	0.54
35:BA:632:A:N3	35:BA:2403:C:H1'	2.22	0.54
35:BA:692:C:N3	35:BA:771:G:C2	2.75	0.54
35:BA:991:C:H42	35:BA:1163:G:H1	1.54	0.54
35:BA:1490:A:C2	38:BD:75:ILE:HD12	2.41	0.54
40:BF:143:ALA:O	40:BF:146:ALA:HB3	2.07	0.54
41:BG:4:ASP:HA	41:BG:8:LYS:CD	2.34	0.54
42:BH:149:ARG:CG	42:BH:162:ILE:HD11	2.37	0.54
45:BO:40:VAL:HA	45:BO:58:VAL:O	2.07	0.54
46:BP:122:PRO:HB3	46:BP:141:ALA:CB	2.36	0.54
49:BS:65:VAL:O	49:BS:69:VAL:HG12	2.07	0.54
50:BT:63:VAL:O	50:BT:73:GLU:HA	2.07	0.54
51:BU:92:ARG:C	51:BU:94:ASN:N	2.60	0.54
52:BV:15:GLU:HB3	52:BV:16:PRO:CD	2.36	0.54
52:BV:34:GLU:HB3	52:BV:62:LEU:CD1	2.34	0.54
53:BW:59:VAL:HG12	53:BW:60:ASN:N	2.22	0.54
54:BX:52:VAL:H	54:BX:80:ILE:HG22	1.72	0.54
54:BX:76:ARG:C	54:BX:76:ARG:HD3	2.27	0.54
1:CA:17:U:O4'	1:CA:1080:A:H1'	2.07	0.54
1:CA:955:U:C1'	1:CA:1227:A:H61	2.15	0.54
1:CA:358:U:H2'	1:CA:359:U:C6	2.42	0.54
1:CA:666:G:H1'	1:CA:741:G:N2	2.22	0.54
3:CC:83:ARG:C	3:CC:87:LEU:HG	2.26	0.54
5:CE:147:ASP:CA	5:CE:150:ARG:HB3	2.37	0.54
6:CF:58:GLY:O	6:CF:60:PHE:CD1	2.60	0.54
8:CH:53:VAL:O	8:CH:54:ASP:HB2	2.07	0.54
8:CH:91:ARG:CG	8:CH:91:ARG:HH11	2.20	0.54
9:CI:16:ARG:O	9:CI:63:ILE:HG22	2.06	0.54
11:CK:92:GLU:O	11:CK:95:ILE:HG12	2.07	0.54
12:CL:34:ARG:HB3	12:CL:61:THR:CG2	2.36	0.54
20:CT:36:LEU:HD22	20:CT:36:LEU:N	2.18	0.54
26:D0:24:LYS:NZ	35:DA:2355:C:O2'	2.41	0.54
29:D3:50:VAL:O	29:D3:51:ALA:C	2.46	0.54
34:D8:11:LYS:HG2	34:D8:11:LYS:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1322:A:O2'	35:DA:1323:U:H5'	2.07	0.54
35:DA:1528(A):A:C2'	35:DA:1529:G:H5''	2.37	0.54
35:DA:15:G:H2'	35:DA:16:G:C8	2.34	0.54
35:DA:1614:A:N1	53:DW:87:PRO:HB3	2.21	0.54
35:DA:1721:G:H2'	35:DA:1722:A:H2'	1.88	0.54
35:DA:1789:A:H2'	35:DA:1790:C:C6	2.38	0.54
35:DA:2591:C:H2'	35:DA:2592:G:H8	1.65	0.54
35:DA:2718:G:H2'	35:DA:2719:G:C8	2.42	0.54
35:DA:2834:G:H5'	35:DA:2835:A:OP2	2.07	0.54
35:DA:409:C:O2'	35:DA:410:G:H5'	2.07	0.54
35:DA:920:G:H2'	35:DA:921:G:C8	2.40	0.54
35:DA:94:C:H5'	35:DA:94(A):G:OP2	2.07	0.54
36:DB:40:U:N3	36:DB:43:C:H5''	2.21	0.54
37:DC:170:ALA:C	37:DC:172:HIS:H	2.11	0.54
38:DD:166:GLN:N	38:DD:166:GLN:NE2	2.54	0.54
39:DE:44:TYR:O	39:DE:45:THR:CB	2.55	0.54
39:DE:47:VAL:HG23	39:DE:84:PHE:O	2.07	0.54
39:DE:69:LYS:N	39:DE:69:LYS:HE2	2.22	0.54
40:DF:32:LEU:O	40:DF:33:LEU:C	2.45	0.54
40:DF:51:THR:OG1	40:DF:91:GLY:HA3	2.07	0.54
41:DG:105:LYS:HZ2	41:DG:105:LYS:CB	2.20	0.54
41:DG:57:ALA:HB2	41:DG:90:LEU:CD2	2.33	0.54
41:DG:80:PHE:O	41:DG:81:LYS:HB2	2.05	0.54
44:DN:10:GLU:CD	44:DN:11:PRO:HD2	2.27	0.54
44:DN:16:ILE:HG23	44:DN:54:VAL:CG2	2.36	0.54
44:DN:77:GLY:O	44:DN:78:TYR:HB3	2.07	0.54
46:DP:23:PRO:O	46:DP:33:ARG:HG2	2.06	0.54
46:DP:51:PHE:O	46:DP:52:GLU:HB2	2.07	0.54
46:DP:95:VAL:O	46:DP:95:VAL:HG23	2.07	0.54
47:DQ:42:ILE:HD13	47:DQ:97:VAL:HG21	1.89	0.54
51:DU:45:TYR:O	51:DU:46:ALA:C	2.45	0.54
55:DY:86:ARG:HB3	55:DY:88:LYS:NZ	2.20	0.54
56:DZ:26:GLY:C	56:DZ:37:VAL:H	2.09	0.54
1:AA:102:G:H2'	1:AA:103:C:H6	1.71	0.54
1:AA:920:U:C1'	1:AA:1080:A:C2	2.89	0.54
1:AA:1417:G:H2'	1:AA:1482:G:N2	2.23	0.54
1:AA:659:U:H2'	1:AA:660:G:H8	1.72	0.54
1:AA:688:G:O2'	1:AA:689:C:H5'	2.06	0.54
1:AA:76:C:H42	1:AA:93:G:H1	1.56	0.54
2:AB:88:ALA:HB2	2:AB:223:ILE:HD11	1.89	0.54
2:AB:82:ARG:HG2	2:AB:82:ARG:HH11	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:170:VAL:HG22	4:AD:171:GLY:H	1.71	0.54
5:AE:43:LEU:CD2	5:AE:132:ALA:HB1	2.31	0.54
5:AE:147:ASP:CA	5:AE:150:ARG:HH11	2.14	0.54
10:AJ:92:THR:HG23	10:AJ:93:GLY:N	2.22	0.54
11:AK:87:THR:HA	11:AK:91:ARG:CG	2.36	0.54
11:AK:65:ALA:HB1	11:AK:98:LEU:CD2	2.37	0.54
12:AL:84:LEU:HB3	12:AL:101:VAL:CG2	2.37	0.54
16:AP:40:ASP:HB3	16:AP:48:TRP:HB3	1.89	0.54
27:B1:47:GLN:NE2	27:B1:64:ALA:CB	2.70	0.54
32:B6:20:ASN:HD22	32:B6:21:TYR:H	1.52	0.54
35:BA:1227:G:OP1	51:BU:13:LYS:HE2	2.07	0.54
35:BA:1596:A:H5'	35:BA:1597:A:OP2	2.07	0.54
35:BA:1748:G:H8	35:BA:1748:G:H5'	1.72	0.54
35:BA:1876:A:H2'	35:BA:1877:A:H8	1.68	0.54
35:BA:202:U:H2'	35:BA:203:C:C6	2.43	0.54
35:BA:2834:G:H5'	35:BA:2835:A:OP2	2.07	0.54
35:BA:37:C:H2'	35:BA:38:A:H8	1.73	0.54
35:BA:469:G:O2'	35:BA:470:A:H5''	2.08	0.54
35:BA:785:G:H2'	35:BA:786:C:C6	2.43	0.54
36:BB:51:G:H2'	36:BB:52:A:O4'	2.07	0.54
38:BD:145:VAL:HB	38:BD:155:LEU:CB	2.38	0.54
38:BD:9:TYR:CD2	38:BD:10:THR:HG22	2.43	0.54
39:BE:179:GLU:OE1	39:BE:179:GLU:HA	2.07	0.54
35:BA:2680:C:H5'	39:BE:189:PRO:HA	1.89	0.54
40:BF:39:TRP:CB	40:BF:101:LEU:HD22	2.37	0.54
40:BF:45:ARG:HG2	40:BF:97:TYR:CG	2.42	0.54
42:BH:88:LEU:O	42:BH:89:ILE:HG23	2.07	0.54
45:BO:20:MET:HE3	45:BO:44:LYS:HG3	1.89	0.54
47:BQ:141:GLN:NE2	56:BZ:72:ARG:HG2	2.21	0.54
48:BR:87:TYR:O	48:BR:89:ASP:N	2.40	0.54
50:BT:107:ASP:OD2	50:BT:109:GLU:HG3	2.06	0.54
52:BV:58:VAL:HG12	52:BV:101:GLY:O	2.07	0.54
54:BX:73:ARG:N	54:BX:74:PRO:CD	2.59	0.54
56:BZ:105:VAL:O	56:BZ:141:VAL:HG13	2.06	0.54
1:CA:102:G:H2'	1:CA:103:C:H6	1.71	0.54
1:CA:1116:C:H2'	1:CA:1117:G:C4'	2.37	0.54
1:CA:1298:C:C6	7:CG:114:ARG:NH1	2.75	0.54
1:CA:159:G:N1	1:CA:163:C:N4	2.56	0.54
1:CA:189(F):U:O4	17:CQ:62:SER:HB3	2.06	0.54
1:CA:261:U:C6	20:CT:79:ARG:NH1	2.75	0.54
1:CA:522:C:H2'	1:CA:523:A:C8	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:78:G:H22	1:CA:91:C:H42	1.53	0.54
2:CB:9:GLU:H	2:CB:9:GLU:CD	2.10	0.54
8:CH:124:ALA:O	8:CH:128:GLY:N	2.38	0.54
11:CK:102:GLY:O	11:CK:103:LEU:C	2.46	0.54
13:CM:93:ARG:HE	13:CM:93:ARG:HA	1.71	0.54
18:CR:36:ASN:O	18:CR:39:VAL:HG23	2.07	0.54
18:CR:35:ARG:C	18:CR:37:VAL:N	2.60	0.54
18:CR:47:THR:HA	18:CR:83:GLU:O	2.07	0.54
24:CX:17:U:O2'	24:CX:18:C:H5'	2.06	0.54
27:D1:76:ARG:HA	27:D1:76:ARG:NE	2.20	0.54
29:D3:4:LEU:HD21	29:D3:56:VAL:HG13	1.88	0.54
31:D5:29:THR:O	31:D5:42:PRO:HD3	2.07	0.54
35:DA:1152:C:H1'	51:DU:77:SER:HB3	1.87	0.54
35:DA:1417:C:C2'	35:DA:1418:G:H5'	2.37	0.54
35:DA:1567:A:H5'	38:DD:58:HIS:CD2	2.42	0.54
35:DA:2512:C:H2'	35:DA:2513:G:O4'	2.07	0.54
35:DA:658:C:H2'	35:DA:659:C:H6	1.73	0.54
35:DA:677:A:N1	35:DA:802:A:C5	2.75	0.54
36:DB:55:U:O2'	36:DB:56:G:H5'	2.07	0.54
38:DD:203:ASN:O	38:DD:204:ILE:O	2.24	0.54
40:DF:68:LYS:O	40:DF:69:HIS:HB2	2.07	0.54
40:DF:89:VAL:C	40:DF:91:GLY:H	2.10	0.54
42:DH:128:PRO:HG2	42:DH:129:THR:CG2	2.32	0.54
43:DI:13:GLY:O	43:DI:15:VAL:N	2.40	0.54
43:DI:33:ARG:O	43:DI:35:LEU:HG	2.08	0.54
46:DP:64:LYS:C	46:DP:66:GLY:N	2.61	0.54
50:DT:106:SER:O	50:DT:107:ASP:HB3	2.07	0.54
50:DT:27:THR:O	50:DT:28:VAL:CB	2.55	0.54
51:DU:51:LYS:HA	51:DU:51:LYS:HE2	1.88	0.54
54:DX:51:VAL:HG13	54:DX:80:ILE:N	2.22	0.54
54:DX:65:ARG:CA	54:DX:65:ARG:NE	2.69	0.54
55:DY:28:LYS:HA	55:DY:39:VAL:N	2.23	0.54
55:DY:7:VAL:CG2	55:DY:8:LYS:HD2	2.37	0.54
1:AA:1424:C:H2'	1:AA:1425:U:C6	2.40	0.54
1:AA:1525:G:O2'	1:AA:1526:G:H5'	2.07	0.54
1:AA:552:U:H4'	12:AL:87:GLY:H	1.72	0.54
1:AA:1096:C:H5''	2:AB:137:ARG:NH2	2.23	0.54
2:AB:89:GLY:O	2:AB:154:LEU:HD13	2.07	0.54
3:AC:56:ASP:O	3:AC:57:ILE:HG13	2.08	0.54
4:AD:15:GLU:C	4:AD:17:VAL:H	2.08	0.54
4:AD:177:ASP:O	4:AD:180:GLY:N	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:58:GLY:O	6:AF:60:PHE:CD1	2.60	0.54
13:AM:108:ARG:H	13:AM:108:ARG:CD	2.12	0.54
20:AT:25:ARG:HH11	20:AT:25:ARG:HG3	1.73	0.54
27:B1:14:VAL:HG22	27:B1:15:ALA:N	2.21	0.54
28:B2:50:ILE:HG23	28:B2:54:LYS:CD	2.32	0.54
29:B3:51:ALA:O	29:B3:53:LEU:N	2.39	0.54
31:B5:12:SER:O	31:B5:13:LYS:C	2.45	0.54
35:BA:120:U:H1'	35:BA:149:A:N7	2.22	0.54
35:BA:1272:A:OP2	35:BA:1647:G:OP1	2.25	0.54
35:BA:1684:C:H42	35:BA:1704:G:H1	1.55	0.54
35:BA:1685:C:H2'	35:BA:1686:C:H6	1.70	0.54
35:BA:1884:A:C3'	35:BA:1885:A:H5''	2.37	0.54
35:BA:2097:C:O2'	35:BA:2098:U:H5'	2.08	0.54
35:BA:2462:U:O2'	35:BA:2463:C:H5'	2.07	0.54
35:BA:2846:G:H2'	35:BA:2847:U:H6	1.72	0.54
35:BA:2893:G:C5'	35:BA:2894:G:H5'	2.16	0.54
35:BA:532:A:N3	35:BA:532:A:H2'	2.21	0.54
35:BA:549:G:C3'	35:BA:551:G:H5''	2.38	0.54
35:BA:549:G:H2'	35:BA:551:G:C4'	2.37	0.54
35:BA:985:C:H2'	35:BA:985:C:O2	2.08	0.54
35:BA:744:G:OP1	39:BE:132:HIS:HB3	2.07	0.54
39:BE:38:THR:HG23	39:BE:39:PRO:HD2	1.88	0.54
40:BF:84:VAL:O	40:BF:85:GLY:C	2.46	0.54
43:BI:49:ALA:HA	43:BI:52:ARG:CG	2.38	0.54
44:BN:10:GLU:CD	44:BN:11:PRO:HD2	2.27	0.54
46:BP:23:PRO:O	46:BP:33:ARG:HG2	2.07	0.54
48:BR:13:HIS:O	48:BR:14:SER:O	2.26	0.54
48:BR:73:VAL:HG23	48:BR:74:LYS:CD	2.38	0.54
49:BS:69:VAL:HG13	49:BS:70:GLY:N	2.21	0.54
50:BT:38:ASN:ND2	50:BT:40:THR:OG1	2.40	0.54
51:BU:49:HIS:O	51:BU:52:ARG:HB2	2.07	0.54
54:BX:21:PHE:HE1	54:BX:26:TYR:HB3	1.71	0.54
56:BZ:126:VAL:HG12	56:BZ:164:ALA:H	1.72	0.54
1:CA:1053:G:C3'	1:CA:1054:C:H5'	2.38	0.54
1:CA:920:U:C1'	1:CA:1080:A:C2	2.87	0.54
1:CA:1202:G:H2'	1:CA:1203:C:C5'	2.37	0.54
1:CA:1269:A:H2	1:CA:1312:G:N3	2.04	0.54
1:CA:527:G:C2'	1:CA:528:C:H5'	2.37	0.54
1:CA:946:A:H2'	1:CA:947:G:H8	1.70	0.54
1:CA:971:G:H4'	1:CA:972:C:C5'	2.36	0.54
3:CC:35:GLU:O	3:CC:39:ILE:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:76:VAL:HG23	3:CC:77:ILE:N	2.22	0.54
3:CC:86:VAL:O	3:CC:89:GLU:HB3	2.08	0.54
4:CD:108:LEU:HD11	4:CD:174:LEU:HD22	1.89	0.54
5:CE:150:ARG:HB2	5:CE:150:ARG:HH11	1.69	0.54
9:CI:77:ILE:HG23	9:CI:81:ILE:HD11	1.90	0.54
11:CK:27:ASN:OD1	11:CK:28:THR:N	2.41	0.54
13:CM:92:HIS:CE1	13:CM:98:VAL:HG21	2.42	0.54
15:CO:75:PRO:O	15:CO:79:ARG:HG3	2.08	0.54
25:CY:164:ILE:CD1	25:CY:164:ILE:H	2.17	0.54
25:CY:53:ASN:OD1	25:CY:54:GLN:HG3	2.07	0.54
26:D0:19:LYS:C	26:D0:20:ARG:HD3	2.27	0.54
26:D0:29:GLN:HB2	26:D0:67:VAL:CG2	2.37	0.54
35:DA:1202:C:H2'	35:DA:1203:G:H5'	1.88	0.54
35:DA:1907:G:O2'	35:DA:1908:C:H5'	2.07	0.54
35:DA:2248:C:H2'	35:DA:2249:U:H5'	1.90	0.54
35:DA:271(N):U:H5''	35:DA:271(O):C:H5'	1.89	0.54
35:DA:271(Q):G:H2'	35:DA:271(R):G:C8	2.43	0.54
31:D5:29:THR:HG21	35:DA:2814:C:O2'	2.07	0.54
35:DA:2842:G:O2'	35:DA:2843:G:H5'	2.07	0.54
35:DA:363(A):A:N3	35:DA:363(A):A:H2'	2.21	0.54
27:D1:69:LYS:HZ2	35:DA:372:G:P	2.31	0.54
35:DA:635:C:O2'	35:DA:636:G:H5'	2.07	0.54
35:DA:66:C:H2'	35:DA:67:U:H5'	1.89	0.54
35:DA:66:C:H2'	35:DA:67:U:C5'	2.38	0.54
35:DA:70:G:H2'	35:DA:113:G:O2'	2.08	0.54
35:DA:750:A:H2'	35:DA:751:A:H5''	1.90	0.54
35:DA:729:G:HO2'	35:DA:763:G:H4'	1.71	0.54
35:DA:768:G:H2'	35:DA:769:G:H8	1.72	0.54
37:DC:65:PRO:HG2	37:DC:189:ILE:HA	1.89	0.54
39:DE:47:VAL:HG12	39:DE:49:LEU:HD21	1.89	0.54
39:DE:55:ASN:O	39:DE:57:LYS:N	2.40	0.54
40:DF:143:ALA:O	40:DF:146:ALA:HB3	2.07	0.54
35:DA:322:A:OP2	40:DF:169:ASN:HB2	2.07	0.54
44:DN:32:THR:HG22	44:DN:37:LYS:HB3	1.88	0.54
44:DN:65:LYS:HA	44:DN:65:LYS:CE	2.28	0.54
35:DA:943:U:OP1	46:DP:38:GLN:HB3	2.08	0.54
56:DZ:27:VAL:HG11	56:DZ:85:HIS:HE2	1.72	0.54
1:AA:1006:C:O2'	1:AA:1007:C:H5'	2.08	0.54
1:AA:284:G:H2'	1:AA:285:G:C8	2.40	0.54
1:AA:436:C:O2'	1:AA:437:U:P	2.65	0.54
1:AA:491:G:H2'	1:AA:492:G:H8	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:591:U:H2'	1:AA:592:G:H8	1.63	0.54
1:AA:865:A:H5'	1:AA:1078:U:C5	2.43	0.54
1:AA:882:C:H2'	1:AA:883:C:H6	1.71	0.54
2:AB:219:VAL:C	2:AB:222:ILE:HG22	2.28	0.54
2:AB:9:GLU:CD	2:AB:9:GLU:H	2.11	0.54
3:AC:120:VAL:HA	3:AC:123:GLN:NE2	2.22	0.54
7:AG:135:VAL:HG12	7:AG:139:GLU:HG3	1.89	0.54
9:AI:11:LYS:O	9:AI:13:ALA:N	2.41	0.54
9:AI:11:LYS:C	9:AI:13:ALA:H	2.10	0.54
12:AL:119:LYS:C	12:AL:121:GLY:H	2.09	0.54
13:AM:23:TYR:HE1	13:AM:71:ARG:CB	2.20	0.54
3:AC:20:SER:O	14:AN:54:PRO:HG3	2.07	0.54
18:AR:74:ARG:HA	18:AR:79:LEU:HB2	1.89	0.54
19:AS:36:ARG:HH12	19:AS:75:ALA:CB	2.20	0.54
29:B3:41:PRO:HB3	35:BA:852:G:O2'	2.08	0.54
31:B5:16:ARG:HH11	31:B5:16:ARG:CG	2.18	0.54
31:B5:20:ARG:O	31:B5:21:SER:C	2.46	0.54
34:B8:60:LEU:HD23	34:B8:60:LEU:N	2.23	0.54
35:BA:1581:G:H2'	35:BA:1582:C:O4'	2.07	0.54
35:BA:2725:A:O2'	35:BA:2726:U:H2'	2.08	0.54
35:BA:470:A:C2	35:BA:471:A:C4	2.96	0.54
35:BA:1843:C:C1'	38:BD:255:LYS:HZ3	2.20	0.54
40:BF:114:VAL:O	40:BF:117:ARG:N	2.41	0.54
44:BN:42:TRP:N	51:BU:64:ARG:HH12	2.05	0.54
46:BP:64:LYS:C	46:BP:66:GLY:N	2.60	0.54
48:BR:37:THR:HG23	48:BR:40:LYS:HE2	1.90	0.54
1:AA:1442(A):G:H21	50:BT:119:LYS:N	2.05	0.54
51:BU:91:ASP:OD2	51:BU:96:ALA:HA	2.07	0.54
53:BW:34:ASN:HA	53:BW:37:ARG:HB3	1.88	0.54
54:BX:77:LYS:HD3	54:BX:78:LYS:HG3	1.90	0.54
55:BY:43:ASN:O	55:BY:44:ILE:O	2.26	0.54
56:BZ:48:PHE:O	56:BZ:50:GLN:N	2.40	0.54
1:CA:66:G:N2	1:CA:172:A:H2	2.05	0.54
1:CA:751:U:C2'	1:CA:752:G:H5'	2.37	0.54
3:CC:134:ILE:O	3:CC:137:ALA:HB3	2.06	0.54
3:CC:147:LYS:HB3	3:CC:203:PHE:CE2	2.43	0.54
4:CD:129:ASN:HB2	4:CD:131:ARG:NH2	2.22	0.54
9:CI:10:ARG:O	9:CI:11:LYS:HB3	2.07	0.54
9:CI:28:VAL:HG12	9:CI:29:ASN:H	1.69	0.54
12:CL:86:ARG:CG	12:CL:87:GLY:N	2.71	0.54
13:CM:68:GLY:HA2	13:CM:71:ARG:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:26:LEU:HD21	18:CR:42:ARG:CZ	2.38	0.54
19:CS:41:VAL:HB	19:CS:44:MET:CB	2.36	0.54
20:CT:11:SER:HA	20:CT:13:LEU:HD12	1.90	0.54
20:CT:54:LYS:HA	20:CT:57:ARG:NH1	2.22	0.54
29:D3:59:VAL:CG1	29:D3:60:GLU:N	2.70	0.54
35:DA:1014:U:H2'	35:DA:1015:G:C8	2.43	0.54
35:DA:1301:A:HO2'	35:DA:1302:A:P	2.30	0.54
35:DA:1411:C:O2'	35:DA:1412:A:H8	1.90	0.54
35:DA:2061:G:OP1	40:DF:68:LYS:NZ	2.40	0.54
35:DA:2645:G:C3'	35:DA:2646:C:H5'	2.27	0.54
35:DA:2732:G:H3'	35:DA:2733:A:H5'	1.90	0.54
35:DA:2863:C:C3'	35:DA:2864:G:H5''	2.38	0.54
35:DA:380:U:H2'	35:DA:381:G:C8	2.41	0.54
35:DA:422:A:H2'	35:DA:423:A:C8	2.42	0.54
35:DA:481:G:H1'	35:DA:506:G:H21	1.69	0.54
35:DA:668:G:O6	35:DA:670:A:H2'	2.08	0.54
35:DA:893:C:H5	35:DA:894:C:C2	2.24	0.54
38:DD:208:LYS:O	38:DD:210:GLY:O	2.25	0.54
38:DD:264:LYS:HG3	38:DD:265:PRO:HD2	1.89	0.54
41:DG:170:ARG:NH1	41:DG:170:ARG:HG3	2.21	0.54
43:DI:3:VAL:HG12	43:DI:37:VAL:O	2.06	0.54
46:DP:124:LYS:HA	46:DP:143:GLY:CA	2.38	0.54
50:DT:34:VAL:O	50:DT:34:VAL:HG12	2.07	0.54
52:DV:22:VAL:O	52:DV:23:GLU:CB	2.55	0.54
54:DX:77:LYS:HD3	54:DX:78:LYS:HG3	1.89	0.54
1:AA:1523:G:C5	1:AA:1524:C:C5	2.95	0.54
1:AA:328:C:C2'	1:AA:328:C:O2	2.56	0.54
1:AA:328:C:H4'	1:AA:329:A:C5'	2.36	0.54
8:AH:63:LEU:HB3	8:AH:65:TYR:CE1	2.42	0.54
12:AL:60:LEU:H	12:AL:60:LEU:HD22	1.71	0.54
12:AL:60:LEU:HD23	12:AL:64:TYR:O	2.08	0.54
13:AM:22:ILE:HG21	13:AM:25:ILE:HD12	1.90	0.54
14:AN:42:ILE:O	14:AN:43:CYS:C	2.44	0.54
15:AO:39:LEU:HD13	15:AO:39:LEU:O	2.08	0.54
15:AO:33:THR:HG23	15:AO:63:ARG:NH1	2.23	0.54
18:AR:47:THR:HA	18:AR:83:GLU:O	2.07	0.54
22:AV:36:A:N1	24:AX:16:U:C4	2.75	0.54
28:B2:56:GLN:NE2	28:B2:56:GLN:HA	2.20	0.54
35:BA:1885:A:H3'	35:BA:1886:C:H6	1.72	0.54
35:BA:2059:A:H5'	35:BA:2060:A:OP2	2.07	0.54
35:BA:2260:C:O2'	35:BA:2261:C:H5'	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2345:G:H5''	35:BA:2347:C:O4'	2.07	0.54
35:BA:2632:A:H1'	39:BE:61:ARG:HH12	1.72	0.54
35:BA:514:A:H2'	35:BA:515:A:C8	2.42	0.54
38:BD:268:ARG:HH12	38:BD:269:PHE:HE1	1.54	0.54
42:BH:99:VAL:O	42:BH:101:ARG:N	2.40	0.54
42:BH:148:ILE:N	42:BH:148:ILE:HD13	2.22	0.54
42:BH:65:HIS:ND1	42:BH:66:GLY:N	2.56	0.54
43:BI:6:LEU:O	43:BI:15:VAL:HG12	2.08	0.54
44:BN:32:THR:HG23	44:BN:37:LYS:HB3	1.88	0.54
44:BN:42:TRP:CD2	44:BN:44:PRO:HD3	2.42	0.54
44:BN:43:THR:O	44:BN:45:ASN:N	2.40	0.54
45:BO:97:ARG:HH11	45:BO:97:ARG:HG3	1.72	0.54
50:BT:109:GLU:HA	50:BT:112:ARG:CD	2.38	0.54
51:BU:24:TYR:HB2	51:BU:29:SER:HB3	1.90	0.54
51:BU:31:SER:C	51:BU:33:ARG:H	2.09	0.54
1:CA:1043:C:H2'	1:CA:1044:A:H8	1.73	0.54
1:CA:1089:G:O2'	1:CA:1090:U:H5'	2.08	0.54
1:CA:1243:C:OP2	21:CU:10:ARG:CZ	2.55	0.54
1:CA:1253:G:H2'	1:CA:1254:C:H6	1.73	0.54
1:CA:192:U:H4'	20:CT:102:GLY:O	2.07	0.54
1:CA:519:C:O2'	1:CA:520:A:H5'	2.08	0.54
1:CA:927:G:H2'	1:CA:928:G:H8	1.71	0.54
1:CA:972:C:H5'	10:CJ:57:LYS:HZ3	1.73	0.54
3:CC:125:GLU:CG	3:CC:189:ALA:HA	2.38	0.54
4:CD:18:LYS:HE3	4:CD:31:CYS:SG	2.48	0.54
4:CD:30:LYS:HB3	4:CD:35:ARG:HH11	1.71	0.54
6:CF:60:PHE:O	6:CF:61:LEU:HD12	2.06	0.54
8:CH:83:ILE:O	8:CH:83:ILE:HG23	2.07	0.54
10:CJ:22:LYS:NZ	10:CJ:23:ILE:HG12	2.22	0.54
10:CJ:36:GLY:O	10:CJ:72:VAL:HG22	2.06	0.54
12:CL:60:LEU:O	12:CL:62:SER:N	2.39	0.54
17:CQ:29:HIS:CE1	17:CQ:31:LEU:HB3	2.43	0.54
17:CQ:71:PHE:CD2	17:CQ:71:PHE:N	2.76	0.54
18:CR:71:LYS:O	18:CR:75:ILE:HG13	2.08	0.54
19:CS:49:ILE:N	19:CS:49:ILE:HD12	2.22	0.54
32:D6:15:GLU:O	32:D6:16:CYS:SG	2.61	0.54
34:D8:49:VAL:CG1	34:D8:53:PRO:HD3	2.38	0.54
35:DA:1036:G:O2'	35:DA:1037:G:H5'	2.08	0.54
35:DA:132:G:O2'	35:DA:133:C:H5'	2.08	0.54
35:DA:1362:C:O2'	35:DA:1363:C:H5'	2.08	0.54
35:DA:1717:G:C3'	35:DA:1718:G:H5''	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1803:A:H4'	38:DD:259:THR:HG23	1.88	0.54
35:DA:2111:C:H1'	35:DA:2118:U:O4'	2.08	0.54
35:DA:648:G:H4'	35:DA:2351:G:H5''	1.89	0.54
35:DA:251:A:H5''	46:DP:51:PHE:HZ	1.71	0.54
35:DA:271(J):C:C3'	35:DA:271(K):U:H5''	2.38	0.54
35:DA:2821:A:OP2	35:DA:2822:G:OP2	2.25	0.54
35:DA:2852:G:H2'	35:DA:2853:C:H6	1.70	0.54
35:DA:360:G:O2'	35:DA:361:G:H5'	2.08	0.54
35:DA:666:G:H4'	46:DP:49:ARG:HH21	1.71	0.54
35:DA:833:U:H5''	46:DP:48:PRO:HB2	1.90	0.54
36:DB:117:G:C5'	49:DS:55:ALA:HB1	2.38	0.54
37:DC:76:ALA:H	37:DC:94:VAL:HG13	1.73	0.54
38:DD:35:LYS:NZ	38:DD:104:TYR:HB2	2.22	0.54
38:DD:69:ARG:HD3	38:DD:105:ILE:HD12	1.88	0.54
38:DD:260:ARG:HG2	38:DD:260:ARG:HH11	1.73	0.54
38:DD:35:LYS:HE2	38:DD:104:TYR:CG	2.42	0.54
38:DD:9:TYR:HD2	38:DD:10:THR:HG22	1.73	0.54
40:DF:168:ARG:HA	40:DF:175:THR:HG21	1.90	0.54
42:DH:159:GLU:O	42:DH:160:LYS:HG2	2.08	0.54
45:DO:6:THR:CG2	45:DO:7:TYR:H	2.19	0.54
45:DO:87:ILE:CG2	45:DO:88:ASN:N	2.71	0.54
46:DP:65:ARG:HH11	46:DP:65:ARG:HG3	1.72	0.54
52:DV:18:LEU:HA	52:DV:97:LYS:HZ1	1.73	0.54
52:DV:4:ILE:HA	52:DV:12:TYR:O	2.08	0.54
53:DW:17:VAL:O	53:DW:19:LEU:N	2.41	0.54
35:DA:1341:U:O3'	54:DX:55:ASN:HB3	2.07	0.54
55:DY:49:VAL:O	55:DY:53:PRO:HG3	2.07	0.54
1:AA:1131:G:H2'	1:AA:1132:C:C6	2.42	0.54
1:AA:237:C:O2'	1:AA:238:G:H5'	2.07	0.54
1:AA:473:G:OP1	16:AP:81:ARG:HB2	2.07	0.54
1:AA:859:A:H2'	1:AA:860:A:O4'	2.06	0.54
2:AB:144:ARG:HG3	2:AB:145:LEU:N	2.22	0.54
3:AC:118:GLN:O	3:AC:122:GLU:HG3	2.08	0.54
5:AE:39:GLY:C	5:AE:69:VAL:HB	2.28	0.54
6:AF:8:ILE:HD13	6:AF:26:ILE:HD13	1.89	0.54
10:AJ:32:ALA:HB1	10:AJ:75:ILE:HG13	1.89	0.54
10:AJ:33:GLN:O	10:AJ:75:ILE:HG23	2.08	0.54
11:AK:17:GLY:HA3	11:AK:79:SER:O	2.07	0.54
12:AL:90:VAL:C	12:AL:92:ASP:H	2.11	0.54
13:AM:61:GLU:O	13:AM:61:GLU:HG2	2.07	0.54
1:AA:237:C:H5''	17:AQ:25:ARG:NH1	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:100:ASN:H	18:AR:23:LYS:NZ	2.05	0.54
18:AR:32:ARG:HA	18:AR:69:THR:HG21	1.90	0.54
25:AY:78:ALA:HA	25:AY:81:LYS:CG	2.38	0.54
27:B1:10:LYS:HG3	27:B1:11:ARG:H	1.73	0.54
29:B3:27:GLY:O	29:B3:35:ARG:HD2	2.07	0.54
33:B7:34:ARG:HE	33:B7:39:ARG:NE	2.06	0.54
35:BA:1244:G:O2'	35:BA:1245:G:H5'	2.08	0.54
35:BA:1281:G:H1	35:BA:1286:A:N6	2.06	0.54
35:BA:1299:G:H5''	35:BA:1300:U:OP1	2.08	0.54
35:BA:1490:A:H5'	35:BA:1491:G:OP2	2.07	0.54
35:BA:1811:G:O2'	35:BA:1812:A:H5'	2.07	0.54
35:BA:2828:C:H2'	35:BA:2829:C:C6	2.42	0.54
35:BA:668:G:O6	35:BA:670:A:H2'	2.07	0.54
35:BA:952:G:C6	35:BA:953:A:N7	2.76	0.54
37:BC:170:ALA:C	37:BC:172:HIS:H	2.11	0.54
39:BE:116:VAL:CG2	39:BE:117:MET:N	2.71	0.54
39:BE:119:ARG:NH1	39:BE:159:HIS:O	2.41	0.54
39:BE:141:ILE:HG12	39:BE:142:GLY:H	1.72	0.54
39:BE:64:LYS:C	39:BE:66:HIS:N	2.61	0.54
40:BF:184:TYR:CD2	40:BF:185:ASP:N	2.75	0.54
40:BF:9:ILE:HG12	40:BF:14:PRO:C	2.28	0.54
41:BG:150:ASP:O	41:BG:151:ALA:HB2	2.08	0.54
41:BG:172:LEU:HA	41:BG:175:LEU:HD12	1.88	0.54
42:BH:151:ILE:N	42:BH:151:ILE:HD13	2.22	0.54
43:BI:4:ILE:O	43:BI:36:ALA:HB1	2.06	0.54
43:BI:68:LEU:HG	43:BI:72:LEU:HD23	1.90	0.54
44:BN:26:LEU:HG	44:BN:30:ILE:HD11	1.89	0.54
45:BO:114:ILE:HD12	45:BO:114:ILE:N	2.04	0.54
35:BA:1952:A:C5	45:BO:22:ILE:HD12	2.43	0.54
45:BO:34:THR:O	45:BO:35:VAL:C	2.46	0.54
47:BQ:114:ALA:C	47:BQ:116:GLU:N	2.60	0.54
49:BS:27:SER:N	49:BS:38:GLN:O	2.41	0.54
50:BT:109:GLU:CB	50:BT:113:LYS:HE3	2.24	0.54
50:BT:120:ARG:HA	50:BT:123:GLN:HG2	1.90	0.54
50:BT:50:ILE:HD12	50:BT:50:ILE:N	2.21	0.54
50:BT:65:LYS:HA	50:BT:65:LYS:HZ1	1.69	0.54
28:B2:29:LYS:HZ2	54:BX:9:LEU:HA	1.71	0.54
55:BY:76:CYS:HB3	55:BY:96:ILE:HD11	1.89	0.54
1:CA:1402:C:O2'	1:CA:1403:C:H5'	2.08	0.54
1:CA:240:C:H2'	1:CA:241:C:H6	1.73	0.54
1:CA:506:G:H2'	1:CA:507:C:C6	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:574:A:N3	1:CA:883:C:H1'	2.22	0.54
1:CA:968:A:H4'	1:CA:969:A:OP2	2.07	0.54
1:CA:973:G:H1'	10:CJ:54:PHE:CE1	2.43	0.54
2:CB:69:LEU:HB3	2:CB:162:ILE:CG2	2.38	0.54
3:CC:118:GLN:O	3:CC:122:GLU:HG3	2.08	0.54
3:CC:19:GLU:O	3:CC:56:ASP:HA	2.08	0.54
3:CC:182:ILE:HG23	3:CC:203:PHE:N	2.22	0.54
4:CD:5:ILE:CG2	4:CD:6:GLY:H	2.13	0.54
7:CG:25:ALA:HA	7:CG:28:ASN:HD22	1.72	0.54
11:CK:34:ASP:HB3	11:CK:40:ILE:HD11	1.90	0.54
12:CL:84:LEU:HB3	12:CL:101:VAL:CG2	2.38	0.54
12:CL:9:GLN:O	12:CL:12:ARG:N	2.39	0.54
13:CM:3:ARG:HA	13:CM:9:ILE:CG1	2.37	0.54
16:CP:48:TRP:CE3	16:CP:49:LEU:HB3	2.42	0.54
18:CR:30:ASP:C	18:CR:32:ARG:H	2.11	0.54
29:D3:32:GLN:HB2	35:DA:1158:C:H4'	1.90	0.54
35:DA:1452:A:O2'	35:DA:1453:U:H2'	2.08	0.54
35:DA:2462:U:O2'	35:DA:2463:C:H5'	2.08	0.54
35:DA:2569:G:O2'	35:DA:2570:G:H5'	2.07	0.54
35:DA:2854:G:H2'	35:DA:2855:C:H6	1.68	0.54
35:DA:58:G:H1	35:DA:69:C:H42	1.55	0.54
36:DB:56:G:H4'	36:DB:57:A:C8	2.43	0.54
38:DD:133:LEU:O	38:DD:134:ARG:C	2.46	0.54
39:DE:131:ALA:HB1	39:DE:134:ILE:HD11	1.89	0.54
35:DA:2578:G:H1'	39:DE:139:GLY:O	2.08	0.54
40:DF:20:LEU:HD12	40:DF:199:TRP:CH2	2.42	0.54
45:DO:13:ASN:HD22	45:DO:97:ARG:CB	2.20	0.54
50:DT:107:ASP:OD2	50:DT:109:GLU:HG3	2.07	0.54
50:DT:28:VAL:HG11	50:DT:46:GLU:OE1	2.07	0.54
51:DU:26:GLY:C	51:DU:28:ARG:N	2.60	0.54
51:DU:65:ILE:HD12	51:DU:65:ILE:H	1.73	0.54
56:DZ:10:ARG:HB2	56:DZ:37:VAL:HA	1.90	0.54
1:AA:1155:G:C2'	1:AA:1156:G:H5'	2.38	0.54
1:AA:1170:A:H2'	1:AA:1171:G:H5'	1.90	0.54
1:AA:1220:G:O2'	1:AA:1221:G:H5'	2.08	0.54
1:AA:1320:C:H2'	1:AA:1321:C:O4'	2.08	0.54
1:AA:38:G:C2	1:AA:397:A:C2	2.96	0.54
1:AA:428:G:O4'	1:AA:430:A:C8	2.61	0.54
1:AA:451:A:N6	1:AA:480:U:H2'	2.23	0.54
1:AA:611:A:O2'	1:AA:612:C:H5'	2.08	0.54
2:AB:175:ARG:O	2:AB:176:GLU:C	2.45	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:35:GLU:O	3:AC:39:ILE:HG13	2.06	0.54
4:AD:104:VAL:O	4:AD:108:LEU:HD13	2.08	0.54
4:AD:105:VAL:CG2	4:AD:126:ILE:HD13	2.36	0.54
5:AE:150:ARG:NH1	5:AE:150:ARG:CB	2.68	0.54
5:AE:80:ILE:HD11	5:AE:91:LEU:HD22	1.89	0.54
8:AH:45:ILE:HA	8:AH:64:LYS:CB	2.37	0.54
10:AJ:16:LEU:HD13	10:AJ:16:LEU:O	2.07	0.54
11:AK:65:ALA:HB1	11:AK:98:LEU:HD23	1.90	0.54
13:AM:23:TYR:OH	13:AM:71:ARG:HD3	2.07	0.54
14:AN:39:LEU:HD11	14:AN:47:LEU:HD12	1.89	0.54
16:AP:80:PHE:HD1	16:AP:80:PHE:H	1.56	0.54
27:B1:34:THR:HG21	35:BA:388:G:P	2.48	0.54
28:B2:57:ILE:O	28:B2:57:ILE:CG1	2.56	0.54
29:B3:3:ARG:HA	29:B3:38:GLU:HA	1.88	0.54
35:BA:1184:G:O2'	35:BA:1185:C:H5'	2.08	0.54
35:BA:1396:U:H2'	35:BA:1396:U:O2	2.07	0.54
35:BA:1469:A:H2'	35:BA:1470:G:O4'	2.06	0.54
35:BA:1778:U:H2'	35:BA:1779:U:C6	2.43	0.54
35:BA:260:G:N2	35:BA:261:G:H1'	2.23	0.54
35:BA:271(G):C:O2'	35:BA:271(H):G:H5'	2.08	0.54
35:BA:354:G:H8	35:BA:354:G:O5'	1.90	0.54
35:BA:534:U:H2'	35:BA:535:C:C6	2.43	0.54
35:BA:623:G:H2'	35:BA:624:C:C6	2.41	0.54
35:BA:692:C:H2'	35:BA:693:C:H6	1.71	0.54
35:BA:88:G:N3	35:BA:88:G:H2'	2.23	0.54
35:BA:2578:G:H1'	39:BE:139:GLY:O	2.08	0.54
39:BE:55:ASN:O	39:BE:57:LYS:N	2.41	0.54
46:BP:132:LYS:O	46:BP:136:GLU:HG2	2.07	0.54
47:BQ:20:ALA:CB	47:BQ:99:PRO:O	2.56	0.54
48:BR:55:ALA:HB1	48:BR:84:ALA:HB2	1.87	0.54
48:BR:74:LYS:HA	48:BR:77:ARG:HD2	1.90	0.54
52:BV:2:PHE:HB2	52:BV:42:GLY:O	2.07	0.54
55:BY:49:VAL:O	55:BY:50:ARG:HB2	2.08	0.54
55:BY:49:VAL:O	55:BY:53:PRO:HG3	2.07	0.54
1:CA:831:U:H2'	1:CA:832:C:H6	1.72	0.54
1:CA:980:C:H5'	1:CA:981:U:C5	2.43	0.54
3:CC:83:ARG:O	3:CC:86:VAL:HG22	2.08	0.54
5:CE:105:VAL:HB	5:CE:106:PRO:CD	2.38	0.54
5:CE:35:GLY:HA2	5:CE:41:VAL:HG12	1.87	0.54
7:CG:50:ILE:CB	7:CG:58:PRO:HD3	2.34	0.54
8:CH:28:ALA:HA	8:CH:59:LEU:HG	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CI:73:GLN:O	9:CI:77:ILE:HG13	2.08	0.54
10:CJ:7:LYS:HG2	10:CJ:71:LEU:HD13	1.89	0.54
14:CN:37:PHE:HE1	14:CN:53:LEU:HD22	1.73	0.54
17:CQ:48:GLU:O	17:CQ:49:GLU:C	2.46	0.54
1:CA:280:C:C4	17:CQ:91:ARG:NH2	2.75	0.54
18:CR:75:ILE:C	18:CR:76:LEU:HD22	2.28	0.54
19:CS:9:VAL:O	19:CS:9:VAL:HG12	2.08	0.54
25:CY:57:THR:HG22	25:CY:59:THR:HG23	1.89	0.54
27:D1:51:VAL:O	27:D1:60:PHE:N	2.41	0.54
27:D1:83:GLU:HG3	27:D1:86:SER:HB2	1.90	0.54
35:DA:1234:U:O2	35:DA:1234:U:H2'	2.07	0.54
35:DA:1396:U:H2'	35:DA:1396:U:O2	2.07	0.54
35:DA:1405:U:H2'	35:DA:1406:U:H6	1.71	0.54
35:DA:1685:C:H2'	35:DA:1686:C:H6	1.73	0.54
35:DA:1820:U:C2	38:DD:202:LYS:HB3	2.43	0.54
35:DA:1884:A:C3'	35:DA:1885:A:H5''	2.37	0.54
35:DA:2464:C:O2'	35:DA:2465:C:H6	1.90	0.54
35:DA:52:A:O2'	35:DA:53:A:H5'	2.08	0.54
36:DB:7:G:H4'	49:DS:29:PHE:HE2	1.71	0.54
37:DC:22:ILE:CG2	37:DC:25:ALA:HB2	2.38	0.54
38:DD:9:TYR:O	38:DD:10:THR:HG22	2.07	0.54
39:DE:77:ILE:HG22	39:DE:79:ARG:HD2	1.89	0.54
42:DH:155:SER:OG	42:DH:156:ALA:N	2.40	0.54
42:DH:157:TYR:HD1	42:DH:170:ARG:O	1.90	0.54
43:DI:6:LEU:O	43:DI:7:GLU:C	2.46	0.54
45:DO:34:THR:O	45:DO:35:VAL:C	2.46	0.54
48:DR:37:THR:OG1	48:DR:40:LYS:HG3	2.07	0.54
49:DS:69:VAL:HG13	49:DS:70:GLY:N	2.22	0.54
35:DA:583:G:OP2	51:DU:10:ARG:NH1	2.40	0.54
52:DV:2:PHE:HB3	52:DV:42:GLY:CA	2.30	0.54
52:DV:5:VAL:CG2	52:DV:36:PRO:HB2	2.35	0.54
54:DX:7:VAL:HA	54:DX:31:HIS:HB2	1.89	0.54
54:DX:25:LYS:NZ	54:DX:87:GLN:O	2.40	0.54
1:AA:1043:C:H2'	1:AA:1044:A:C8	2.41	0.54
1:AA:1276:G:H2'	1:AA:1277:C:H5'	1.89	0.54
1:AA:272:C:O2'	1:AA:273:A:H5'	2.08	0.54
1:AA:309:G:O2'	1:AA:310:G:H5'	2.08	0.54
1:AA:511:C:C2	1:AA:512:U:C5	2.95	0.54
1:AA:545:C:O2'	1:AA:546:G:H5'	2.08	0.54
1:AA:706:A:C5	1:AA:707:C:C5	2.94	0.54
1:AA:831:U:H2'	1:AA:832:C:H6	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:150:GLU:HA	4:AD:153:ARG:CD	2.38	0.54
4:AD:65:ARG:NH1	4:AD:72:GLU:N	2.56	0.54
5:AE:129:ILE:O	5:AE:132:ALA:N	2.40	0.54
5:AE:15:ARG:HG2	5:AE:26:PHE:CD2	2.42	0.54
1:AA:939:G:H5''	7:AG:102:ARG:CZ	2.36	0.54
7:AG:20:ASP:HB3	7:AG:23:VAL:HG23	1.89	0.54
10:AJ:7:LYS:HG2	10:AJ:71:LEU:HD13	1.89	0.54
20:AT:36:LEU:HD22	20:AT:36:LEU:N	2.18	0.54
25:AY:127:VAL:HA	25:AY:130:ARG:HB2	1.89	0.54
29:B3:50:VAL:O	29:B3:51:ALA:C	2.46	0.54
32:B6:9:LEU:HD23	32:B6:10:LEU:N	2.23	0.54
35:BA:1174:A:H5''	35:BA:1175:U:H5''	1.89	0.54
35:BA:1451:C:H4'	35:BA:1452:A:C8	2.43	0.54
35:BA:1744:C:C2'	35:BA:1745:C:H5'	2.38	0.54
35:BA:174:C:O2	35:BA:174:C:H2'	2.08	0.54
35:BA:1918:A:O2'	35:BA:1920:C:N4	2.41	0.54
35:BA:2020:A:N1	35:BA:2034:U:O4	2.41	0.54
35:BA:2512:C:H4'	39:BE:122:PHE:CE2	2.43	0.54
35:BA:2536:G:C6	35:BA:2537:U:N3	2.76	0.54
35:BA:271(D):G:H1	35:BA:271(T):C:N4	1.80	0.54
35:BA:2780:G:OP1	44:BN:118:LYS:HE2	2.07	0.54
35:BA:2870:C:H2'	35:BA:2871:C:H5'	1.90	0.54
35:BA:575:A:O2'	35:BA:576:U:H5'	2.08	0.54
35:BA:7:G:H4'	44:BN:13:TRP:HH2	1.67	0.54
35:BA:860:U:O2'	35:BA:861:A:H5'	2.08	0.54
35:BA:1828:G:O6	38:BD:222:ARG:HD3	2.08	0.54
38:BD:229:VAL:HG23	38:BD:230:ASP:N	2.21	0.54
35:BA:2682:U:C2	39:BE:22:PRO:HB3	2.43	0.54
40:BF:197:ASP:OD1	40:BF:198:ALA:N	2.41	0.54
40:BF:20:LEU:HD12	40:BF:199:TRP:CH2	2.42	0.54
40:BF:30:PRO:O	40:BF:33:LEU:HB3	2.08	0.54
41:BG:31:VAL:HG22	41:BG:32:PRO:HD2	1.90	0.54
41:BG:41:GLN:HE21	41:BG:153:ARG:HB3	1.72	0.54
41:BG:76:SER:HB3	41:BG:84:LYS:HD2	1.89	0.54
42:BH:159:GLU:O	42:BH:160:LYS:HG2	2.08	0.54
43:BI:15:VAL:O	43:BI:16:GLY:C	2.45	0.54
46:BP:95:VAL:HG23	46:BP:95:VAL:O	2.08	0.54
35:BA:869:G:H1'	47:BQ:8:LYS:HZ2	1.70	0.54
49:BS:40:ILE:HG23	49:BS:46:VAL:O	2.08	0.54
50:BT:62:THR:HB	50:BT:74:ARG:O	2.07	0.54
51:BU:51:LYS:HA	51:BU:51:LYS:HE2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:BZ:70:LEU:O	56:BZ:88:PHE:HD2	1.91	0.54
1:CA:115:G:O2'	1:CA:116:A:OP2	2.23	0.54
1:CA:1413:A:O2'	1:CA:1414:U:H5'	2.07	0.54
1:CA:1521:G:C6	1:CA:1522:U:C4	2.96	0.54
1:CA:189(H):G:H2'	1:CA:189(I):G:C8	2.40	0.54
1:CA:191:G:C4	20:CT:105:SER:HB3	2.42	0.54
1:CA:512:U:H2'	1:CA:513:C:C6	2.43	0.54
1:CA:865:A:H2'	1:CA:866:C:C6	2.43	0.54
3:CC:171:GLY:O	3:CC:172:ARG:O	2.26	0.54
4:CD:80:GLU:HB3	4:CD:84:LYS:HZ1	1.73	0.54
6:CF:100:ASN:O	6:CF:101:ALA:O	2.25	0.54
11:CK:125:PHE:HD1	11:CK:125:PHE:N	2.06	0.54
12:CL:21:LYS:N	12:CL:21:LYS:HD2	2.19	0.54
12:CL:58:VAL:O	12:CL:60:LEU:HD22	2.08	0.54
14:CN:41:ARG:HG3	14:CN:42:ILE:H	1.72	0.54
15:CO:11:VAL:HG13	15:CO:15:PHE:HE1	1.73	0.54
16:CP:28:ARG:NH1	16:CP:28:ARG:HG2	2.22	0.54
25:CY:113:ASP:O	25:CY:116:ARG:HG3	2.08	0.54
25:CY:24:ASN:HB3	25:CY:121:TYR:CD1	2.43	0.54
27:D1:62:VAL:CG2	27:D1:63:ALA:N	2.71	0.54
32:D6:36:LEU:HD13	32:D6:50:ARG:HH12	1.71	0.54
35:DA:2007:C:H2'	35:DA:2008:C:C6	2.41	0.54
35:DA:2037:G:H2'	35:DA:2038:G:H8	1.73	0.54
35:DA:2632:A:H1'	39:DE:61:ARG:HH12	1.73	0.54
35:DA:2643:G:O2'	35:DA:2644:G:H5'	2.08	0.54
35:DA:503:A:C6	35:DA:505:A:C6	2.96	0.54
36:DB:51:G:H2'	36:DB:52:A:O4'	2.07	0.54
38:DD:145:VAL:O	38:DD:153:ALA:HB1	2.08	0.54
40:DF:22:ALA:CA	40:DF:26:ALA:HB2	2.37	0.54
36:DB:42:C:N4	41:DG:91:ARG:NH2	2.55	0.54
46:DP:38:GLN:O	46:DP:39:LYS:HB2	2.08	0.54
50:DT:110:ILE:HA	50:DT:113:LYS:HD2	1.89	0.54
51:DU:107:ALA:O	51:DU:111:GLU:HG2	2.08	0.54
1:AA:1043:C:H2'	1:AA:1044:A:H8	1.71	0.54
1:AA:1242:C:O5'	1:AA:1242:C:H6	1.90	0.54
1:AA:903:G:H2'	1:AA:904:C:C6	2.39	0.54
1:AA:968:A:H4'	1:AA:969:A:OP2	2.07	0.54
6:AF:27:GLN:NE2	6:AF:27:GLN:HA	2.23	0.54
8:AH:65:TYR:HA	8:AH:79:VAL:HG23	1.90	0.54
8:AH:7:ALA:O	8:AH:8:ASP:C	2.45	0.54
10:AJ:34:VAL:HG12	10:AJ:35:SER:N	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:21:ILE:CD1	11:AK:84:VAL:HG12	2.38	0.54
12:AL:32:PHE:HD1	12:AL:86:ARG:HA	1.73	0.54
13:AM:28:ALA:C	13:AM:30:ALA:H	2.09	0.54
13:AM:68:GLY:HA2	13:AM:71:ARG:HB3	1.90	0.54
16:AP:20:VAL:HG22	16:AP:21:VAL:N	2.20	0.54
18:AR:87:ARG:HG2	18:AR:88:LYS:N	2.23	0.54
19:AS:42:PRO:C	19:AS:44:MET:H	2.11	0.54
23:AW:38:A:H2'	23:AW:39:A:C5'	2.38	0.54
25:AY:76:LEU:HD13	25:AY:97:ASP:O	2.08	0.54
28:B2:50:ILE:O	28:B2:51:ARG:CB	2.56	0.54
35:BA:1323:U:H3	35:BA:1331:A:H61	1.54	0.54
35:BA:1637:A:H2'	35:BA:1638:C:C6	2.43	0.54
35:BA:2233:U:H2'	35:BA:2234:G:C8	2.42	0.54
35:BA:2512:C:H2'	35:BA:2513:G:O4'	2.08	0.54
35:BA:2522:U:H2'	35:BA:2523:G:H5''	1.89	0.54
35:BA:2571:C:H5'	35:BA:2572:A:H5'	1.88	0.54
35:BA:292:C:H2'	35:BA:292:C:O2	2.07	0.54
35:BA:671:C:C5	46:BP:36:LYS:NZ	2.75	0.54
38:BD:134:ARG:O	38:BD:136:ILE:N	2.40	0.54
40:BF:125:LEU:HB3	40:BF:196:LEU:HD21	1.89	0.54
40:BF:178:PRO:HG2	40:BF:179:GLU:N	2.16	0.54
41:BG:52:ILE:O	41:BG:54:GLU:N	2.41	0.54
41:BG:85:GLY:C	41:BG:87:PRO:CD	2.77	0.54
42:BH:12:PRO:N	42:BH:15:VAL:HG21	2.23	0.54
43:BI:13:GLY:O	43:BI:15:VAL:N	2.41	0.54
35:BA:8:A:H5''	44:BN:51:PHE:HZ	1.73	0.54
46:BP:48:PRO:O	46:BP:51:PHE:N	2.41	0.54
46:BP:83:VAL:HG23	46:BP:105:LEU:HD22	1.88	0.54
52:BV:82:ARG:NH1	52:BV:84:LYS:HD3	2.22	0.54
55:BY:16:ALA:HA	55:BY:21:LYS:HD2	1.90	0.54
55:BY:83:THR:HG22	55:BY:84:ARG:H	1.73	0.54
1:CA:1405:G:H1'	1:CA:1518:A:HO2'	1.72	0.54
1:CA:385:C:H2'	1:CA:386:C:H6	1.73	0.54
1:CA:732:C:H2'	1:CA:733:A:H5''	1.89	0.54
1:CA:9:G:H2'	1:CA:10:A:C8	2.43	0.54
2:CB:54:THR:HG22	2:CB:58:ILE:CD1	2.36	0.54
2:CB:68:ILE:HG22	2:CB:70:PHE:HD1	1.72	0.54
3:CC:120:VAL:HA	3:CC:123:GLN:NE2	2.23	0.54
4:CD:110:PHE:CD1	4:CD:110:PHE:N	2.75	0.54
4:CD:119:GLN:HG3	4:CD:123:HIS:CD2	2.39	0.54
5:CE:39:GLY:C	5:CE:69:VAL:HB	2.28	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CF:70:ASP:CG	6:CF:71:ARG:H	2.09	0.54
9:CI:105:ASP:C	9:CI:107:ARG:H	2.10	0.54
1:CA:881:G:P	12:CL:12:ARG:NH2	2.81	0.54
13:CM:28:ALA:C	13:CM:30:ALA:H	2.11	0.54
18:CR:25:THR:O	18:CR:25:THR:HG22	2.07	0.54
19:CS:36:ARG:CZ	19:CS:72:GLY:HA2	2.38	0.54
20:CT:26:ASN:HD22	20:CT:27:LYS:H	1.55	0.54
28:D2:12:GLU:O	28:D2:12:GLU:CG	2.56	0.54
33:D7:8:ASN:ND2	33:D7:9:ARG:N	2.52	0.54
35:DA:1225:G:OP1	52:DV:88:ARG:HD2	2.07	0.54
35:DA:1495:A:C8	35:DA:1495:A:OP1	2.61	0.54
35:DA:2079:U:H2'	35:DA:2080:G:C8	2.43	0.54
35:DA:2571:C:O2	35:DA:2571:C:H2'	2.07	0.54
35:DA:265:A:H1'	35:DA:266:G:C1'	2.38	0.54
35:DA:391:G:O2'	35:DA:392:C:H5'	2.08	0.54
35:DA:549:G:C3'	35:DA:551:G:H5''	2.37	0.54
36:DB:30:C:H2'	36:DB:31:C:O4'	2.08	0.54
38:DD:146:GLU:CA	38:DD:153:ALA:HA	2.34	0.54
39:DE:32:PRO:O	39:DE:34:VAL:HG13	2.08	0.54
40:DF:107:LYS:O	40:DF:110:LEU:N	2.41	0.54
35:DA:673:C:H5'	40:DF:54:ARG:NH1	2.22	0.54
40:DF:84:VAL:O	40:DF:85:GLY:C	2.46	0.54
41:DG:20:ILE:CD1	41:DG:25:TYR:HB2	2.38	0.54
42:DH:118:PRO:HG2	42:DH:121:ILE:HD12	1.90	0.54
43:DI:109:ILE:HD13	43:DI:111:PRO:HD3	1.89	0.54
45:DO:68:GLU:N	45:DO:68:GLU:OE2	2.41	0.54
46:DP:140:ALA:O	46:DP:141:ALA:HB3	2.08	0.54
49:DS:93:LYS:HD2	49:DS:93:LYS:O	2.08	0.54
56:DZ:99:TYR:HA	56:DZ:124:ILE:O	2.07	0.54
1:AA:1077:G:N2	1:AA:1079:G:H3'	2.23	0.54
1:AA:1104:G:H2'	1:AA:1105:A:H8	1.72	0.54
1:AA:1116:C:H2'	1:AA:1117:G:C4'	2.38	0.54
1:AA:1208:C:H2'	1:AA:1209:C:C6	2.43	0.54
1:AA:1442(A):G:H4'	1:AA:1442(B):A:O5'	2.08	0.54
1:AA:180:U:H2'	1:AA:181:G:H5''	1.89	0.54
1:AA:601:C:O2'	1:AA:602:A:H5'	2.08	0.54
1:AA:950:U:H2'	1:AA:951:G:H8	1.70	0.54
2:AB:67:THR:CG2	2:AB:155:LEU:HG	2.34	0.54
2:AB:193:ASP:OD2	2:AB:196:LEU:HD21	2.08	0.54
5:AE:72:GLN:O	5:AE:73:ASN:HB3	2.08	0.54
7:AG:148:ASN:C	7:AG:150:ALA:H	2.10	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:27:ILE:HG21	7:AG:40:ALA:HB2	1.90	0.54
7:AG:50:ILE:HG22	7:AG:56:GLN:O	2.07	0.54
25:AY:92:PRO:CA	25:AY:101:ILE:HG23	2.38	0.54
35:BA:1202:C:H2'	35:BA:1203:G:H5'	1.90	0.54
35:BA:1495:A:OP1	35:BA:1495:A:C8	2.61	0.54
35:BA:1543:C:O2	35:BA:1543:C:H2'	2.07	0.54
35:BA:1668:A:N7	35:BA:1674:G:C6	2.76	0.54
35:BA:182:A:H2	35:BA:433:C:H1'	1.73	0.54
35:BA:2286:A:H4'	35:BA:2287:A:O4'	2.07	0.54
35:BA:2410:G:N2	35:BA:2411:A:H1'	2.23	0.54
35:BA:271(U):G:O2'	35:BA:271(V):G:H5'	2.08	0.54
35:BA:514:A:O2'	35:BA:515:A:H5'	2.08	0.54
35:BA:776:G:H4'	35:BA:777:A:O5'	2.08	0.54
36:BB:95:C:O2'	36:BB:96:U:H5'	2.08	0.54
37:BC:77:ILE:O	37:BC:77:ILE:HG12	2.07	0.54
38:BD:257:LEU:HD23	38:BD:258:LYS:N	2.23	0.54
39:BE:57:LYS:HD3	39:BE:59:VAL:HG12	1.90	0.54
39:BE:77:ILE:HG23	39:BE:78:LEU:H	1.72	0.54
39:BE:36:ARG:NH2	39:BE:88:GLY:HA3	2.23	0.54
40:BF:51:THR:OG1	40:BF:91:GLY:HA3	2.07	0.54
42:BH:144:VAL:CG1	42:BH:148:ILE:HD11	2.37	0.54
44:BN:76:SER:O	44:BN:77:GLY:C	2.43	0.54
48:BR:70:LEU:O	48:BR:71:GLN:HB2	2.08	0.54
50:BT:102:ILE:CB	50:BT:110:ILE:HD11	2.38	0.54
50:BT:91:ARG:HB3	50:BT:115:ARG:O	2.09	0.54
52:BV:5:VAL:HG23	52:BV:37:VAL:H	1.73	0.54
53:BW:9:TYR:H	53:BW:102:HIS:CD2	2.26	0.54
56:BZ:118:GLN:N	56:BZ:173:ALA:O	2.40	0.54
56:BZ:29:TYR:O	56:BZ:30:ASN:CB	2.56	0.54
1:CA:423:G:H2'	1:CA:424:G:O4'	2.08	0.54
1:CA:896:C:O2'	1:CA:897:C:H5'	2.08	0.54
2:CB:215:LEU:O	2:CB:219:VAL:HG23	2.08	0.54
2:CB:28:PHE:CD1	2:CB:28:PHE:O	2.61	0.54
3:CC:109:PRO:HG2	3:CC:110:ASN:H	1.73	0.54
3:CC:18:TRP:HE3	3:CC:18:TRP:H	1.53	0.54
4:CD:58:LEU:O	4:CD:59:ARG:C	2.44	0.54
7:CG:145:ALA:O	7:CG:146:GLU:HB2	2.08	0.54
8:CH:4:ASP:OD2	8:CH:7:ALA:N	2.29	0.54
1:CA:963:G:N2	10:CJ:55:LYS:NZ	2.56	0.54
10:CJ:62:HIS:N	10:CJ:62:HIS:CD2	2.74	0.54
11:CK:21:ILE:CD1	11:CK:84:VAL:HG12	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:61:GLU:HG2	13:CM:61:GLU:O	2.08	0.54
16:CP:48:TRP:HE3	16:CP:49:LEU:HB3	1.73	0.54
16:CP:67:THR:O	16:CP:71:ARG:HB2	2.07	0.54
21:CU:18:TYR:CD2	21:CU:24:ARG:HG3	2.43	0.54
25:CY:80:GLU:C	25:CY:82:ALA:H	2.09	0.54
25:CY:76:LEU:HD21	25:CY:99:LEU:CD2	2.35	0.54
27:D1:75:GLU:O	27:D1:76:ARG:NE	2.38	0.54
27:D1:76:ARG:HB3	27:D1:78:LYS:HZ3	1.71	0.54
28:D2:57:ILE:HD11	28:D2:59:ARG:NH1	2.23	0.54
33:D7:30:VAL:HG12	33:D7:33:ARG:HH12	1.73	0.54
35:DA:1244:G:O2'	35:DA:1245:G:H5'	2.07	0.54
33:D7:19:ARG:HD3	35:DA:125:G:H5''	1.89	0.54
35:DA:1639:U:H2'	35:DA:1640:C:C5'	2.38	0.54
35:DA:221:A:O2'	35:DA:222:A:OP2	2.25	0.54
35:DA:2078:C:H1'	35:DA:2434:A:N3	2.22	0.54
35:DA:271(J):C:C2'	35:DA:271(K):U:H5''	2.38	0.54
35:DA:272(C):G:H2'	35:DA:272(D):G:C8	2.41	0.54
35:DA:2801:A:H4'	35:DA:2801(A):A:O5'	2.07	0.54
35:DA:336:C:H4'	55:DY:7:VAL:CG1	2.38	0.54
35:DA:402:A:O2'	35:DA:403:U:H5'	2.08	0.54
38:DD:173:VAL:CG1	38:DD:185:VAL:O	2.56	0.54
35:DA:607:U:OP1	40:DF:102:PRO:HA	2.08	0.54
40:DF:110:LEU:HD21	40:DF:181:LEU:CD2	2.38	0.54
42:DH:99:VAL:O	42:DH:101:ARG:N	2.41	0.54
42:DH:164:TYR:O	42:DH:165:ALA:HB2	2.08	0.54
42:DH:55:PRO:HG2	42:DH:61:HIS:NE2	2.21	0.54
44:DN:26:LEU:CD1	44:DN:30:ILE:HD11	2.37	0.54
44:DN:96:GLU:O	44:DN:97:ARG:C	2.45	0.54
47:DQ:114:ALA:C	47:DQ:116:GLU:N	2.60	0.54
48:DR:53:HIS:HA	48:DR:56:LYS:HB2	1.88	0.54
49:DS:12:PHE:O	49:DS:12:PHE:CG	2.61	0.54
50:DT:110:ILE:HG22	50:DT:111:ARG:N	2.23	0.54
52:DV:1:MET:CE	52:DV:45:THR:H	2.20	0.54
52:DV:71:LEU:O	52:DV:90:PRO:HA	2.07	0.54
1:AA:1468:A:H2'	1:AA:1469:G:O4'	2.07	0.53
1:AA:355:C:O2'	1:AA:356:A:H5'	2.07	0.53
1:AA:444:C:H2'	1:AA:445:G:C8	2.41	0.53
1:AA:674:G:H2'	1:AA:675:A:C8	2.37	0.53
1:AA:676:A:O2'	1:AA:677:U:H5'	2.08	0.53
2:AB:164:VAL:O	2:AB:186:ALA:HB1	2.08	0.53
2:AB:187:LEU:HA	2:AB:201:ILE:O	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:82:ARG:HB2	2:AB:94:ASN:ND2	2.23	0.53
4:AD:110:PHE:N	4:AD:110:PHE:CD1	2.76	0.53
4:AD:33:MET:HA	4:AD:33:MET:CE	2.39	0.53
7:AG:24:THR:O	7:AG:25:ALA:C	2.44	0.53
9:AI:28:VAL:HG12	9:AI:29:ASN:H	1.73	0.53
9:AI:16:ARG:O	9:AI:63:ILE:CG2	2.56	0.53
12:AL:119:LYS:HD3	12:AL:120:TYR:HE1	1.73	0.53
10:AJ:50:ILE:CD1	14:AN:41:ARG:HD2	2.33	0.53
14:AN:44:LEU:HD12	14:AN:44:LEU:O	2.08	0.53
15:AO:6:GLU:N	15:AO:6:GLU:OE1	2.32	0.53
22:AV:27:G:H2'	22:AV:28:G:C8	2.43	0.53
25:AY:26:ALA:C	25:AY:28:LEU:H	2.11	0.53
25:AY:84:ARG:O	25:AY:86:SER:N	2.40	0.53
33:B7:34:ARG:HE	33:B7:39:ARG:HE	1.54	0.53
33:B7:5:TRP:HE1	33:B7:7:PRO:HG3	1.72	0.53
34:B8:26:LYS:NZ	34:B8:47:LYS:HD3	2.23	0.53
34:B8:29:LYS:O	34:B8:30:ARG:C	2.47	0.53
34:B8:53:PRO:C	34:B8:55:ALA:N	2.62	0.53
35:BA:1152:C:H1'	51:BU:77:SER:HB3	1.89	0.53
35:BA:1844:C:C2'	35:BA:1845:G:H5'	2.37	0.53
35:BA:1862:G:O2'	35:BA:1863:G:H5'	2.08	0.53
35:BA:1952:A:C4	45:BO:22:ILE:HD12	2.43	0.53
35:BA:271(N):U:H5''	35:BA:271(O):C:H5'	1.89	0.53
35:BA:2748:A:C2	42:BH:63:SER:HB3	2.43	0.53
35:BA:2773:C:H2'	35:BA:2774:C:C6	2.44	0.53
35:BA:2784:C:H2'	35:BA:2785:C:C6	2.43	0.53
35:BA:338:G:H2'	35:BA:339:U:C6	2.43	0.53
35:BA:493:G:C2'	35:BA:494:G:H5''	2.38	0.53
35:BA:535:C:C2'	35:BA:536:A:H5'	2.39	0.53
38:BD:152:GLY:O	38:BD:154:LYS:HG3	2.08	0.53
38:BD:80:ALA:HB2	38:BD:96:HIS:ND1	2.23	0.53
40:BF:151:SER:C	40:BF:152:GLU:HG3	2.29	0.53
40:BF:39:TRP:HA	40:BF:99:TYR:CE1	2.43	0.53
41:BG:130:ASN:ND2	41:BG:161:THR:O	2.41	0.53
48:BR:37:THR:HA	48:BR:111:LEU:HA	1.90	0.53
52:BV:98:GLU:N	52:BV:98:GLU:OE1	2.40	0.53
53:BW:17:VAL:C	53:BW:19:LEU:N	2.61	0.53
1:CA:1078:U:H2'	1:CA:1079:G:C8	2.43	0.53
1:CA:164:U:O2'	1:CA:165:C:H5'	2.08	0.53
1:CA:528:C:H2'	1:CA:529:G:H8	1.73	0.53
1:CA:527:G:H2'	1:CA:528:C:H5'	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:539:A:H2'	1:CA:540:G:H8	1.73	0.53
1:CA:59:A:H2'	1:CA:59:A:N3	2.22	0.53
1:CA:652:U:H1'	1:CA:653:A:C2	2.42	0.53
7:CG:71:PRO:CG	7:CG:103:TRP:HZ3	2.20	0.53
8:CH:83:ILE:HD13	8:CH:137:VAL:CG2	2.35	0.53
8:CH:97:VAL:HG13	8:CH:98:LYS:N	2.17	0.53
13:CM:64:TRP:NE1	13:CM:66:LEU:HD12	2.23	0.53
13:CM:68:GLY:O	13:CM:71:ARG:N	2.41	0.53
10:CJ:50:ILE:CD1	14:CN:41:ARG:HD2	2.36	0.53
1:CA:1216:G:O3'	14:CN:5:ALA:HB1	2.08	0.53
1:CA:607:A:C4	16:CP:31:LYS:HE3	2.43	0.53
19:CS:4:SER:N	19:CS:6:LYS:HZ1	2.07	0.53
20:CT:41:ILE:C	20:CT:43:LEU:N	2.61	0.53
23:CW:7:G:H5'	23:CW:8:U:H5	1.72	0.53
25:CY:67:VAL:HA	25:CY:99:LEU:O	2.08	0.53
35:DA:1024:G:O5'	35:DA:1024:G:H8	1.91	0.53
35:DA:1441:G:O2'	35:DA:1442:G:H5'	2.08	0.53
35:DA:2128:C:H5'	35:DA:2173:A:C2	2.43	0.53
35:DA:2282:G:C4	35:DA:2425:A:N6	2.76	0.53
35:DA:2784:C:H2'	35:DA:2785:C:C6	2.44	0.53
35:DA:2810:A:H2'	39:DE:61:ARG:NH2	2.23	0.53
35:DA:376:C:O2'	35:DA:377:C:H5'	2.09	0.53
35:DA:909:A:H1'	47:DQ:10:ARG:HH22	1.72	0.53
36:DB:45:A:H2'	36:DB:45:A:N3	2.23	0.53
38:DD:145:VAL:CG1	38:DD:146:GLU:N	2.72	0.53
38:DD:211:ARG:HA	38:DD:214:TRP:CD2	2.43	0.53
38:DD:19:ALA:O	38:DD:21:PHE:CE1	2.61	0.53
38:DD:242:ARG:HG3	38:DD:242:ARG:NH1	2.19	0.53
39:DE:116:VAL:CG2	39:DE:122:PHE:HB2	2.32	0.53
42:DH:102:ALA:HB1	42:DH:115:VAL:O	2.07	0.53
43:DI:38:LEU:H	43:DI:38:LEU:CD1	2.03	0.53
47:DQ:51:ARG:C	47:DQ:54:MET:HB3	2.27	0.53
48:DR:49:ASP:O	48:DR:51:LEU:N	2.41	0.53
50:DT:109:GLU:HA	50:DT:112:ARG:CD	2.37	0.53
50:DT:106:SER:CA	50:DT:110:ILE:HD13	2.33	0.53
50:DT:3:ARG:O	50:DT:4:GLY:C	2.47	0.53
50:DT:57:PHE:O	50:DT:59:THR:HG22	2.08	0.53
50:DT:99:LEU:CD1	50:DT:99:LEU:O	2.56	0.53
51:DU:21:ALA:CB	51:DU:35:ALA:HB1	2.37	0.53
55:DY:44:ILE:CG2	55:DY:45:VAL:N	2.70	0.53
1:AA:129(A):G:H21	1:AA:189(F):U:H5"	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:277:C:O2'	1:AA:278:G:H5'	2.07	0.53
1:AA:314:C:O2'	1:AA:315:A:H5'	2.08	0.53
2:AB:12:GLU:C	2:AB:14:GLY:N	2.61	0.53
4:AD:128:VAL:C	4:AD:130:GLY:N	2.61	0.53
4:AD:91:SER:HA	4:AD:94:LEU:HD12	1.90	0.53
5:AE:126:ARG:O	5:AE:127:ASN:C	2.47	0.53
6:AF:30:LEU:HB3	6:AF:35:ALA:HB3	1.90	0.53
9:AI:50:LEU:O	9:AI:55:ALA:HB3	2.08	0.53
11:AK:23:ALA:HA	11:AK:28:THR:HG23	1.89	0.53
11:AK:65:ALA:HB3	11:AK:97:ALA:CB	2.38	0.53
10:AJ:64:GLU:N	14:AN:59:ALA:HB2	2.23	0.53
16:AP:67:THR:O	16:AP:71:ARG:HB2	2.08	0.53
19:AS:49:ILE:HD12	19:AS:49:ILE:N	2.23	0.53
19:AS:60:VAL:O	19:AS:60:VAL:HG13	2.08	0.53
23:AW:65:G:N2	23:AW:66:C:H1'	2.23	0.53
25:AY:63:PRO:HB3	25:AY:64:ARG:HH22	1.73	0.53
26:B0:27:GLU:HG3	26:B0:68:GLU:HA	1.90	0.53
34:B8:30:ARG:HH21	46:BP:62:LEU:CB	2.17	0.53
35:BA:1188:U:H2'	35:BA:1189:A:H5'	1.90	0.53
35:BA:118:A:N3	35:BA:178:G:H1'	2.24	0.53
35:BA:1419:A:H62	35:BA:1578:U:H3	1.56	0.53
35:BA:1992:G:C6	35:BA:1997:G:N1	2.75	0.53
35:BA:2061:G:OP1	40:BF:68:LYS:NZ	2.41	0.53
35:BA:2065:C:H2'	35:BA:2066:C:C6	2.43	0.53
35:BA:2415:G:C3'	46:BP:66:GLY:HA3	2.38	0.53
35:BA:2679:A:O2'	35:BA:2680:C:H5'	2.07	0.53
35:BA:2845:G:H2'	35:BA:2846:G:H8	1.74	0.53
35:BA:510:C:O2'	35:BA:511:U:H5'	2.09	0.53
35:BA:548:A:O2'	35:BA:549:G:OP1	2.26	0.53
35:BA:614(C):A:N3	40:BF:180:GLY:HA2	2.23	0.53
35:BA:644:A:C2	35:BA:2369:A:H1'	2.42	0.53
38:BD:125:ILE:N	38:BD:125:ILE:CD1	2.71	0.53
38:BD:169:GLU:O	38:BD:171:ASP:N	2.36	0.53
38:BD:260:ARG:NH2	38:BD:264:LYS:HD3	2.23	0.53
39:BE:81:ILE:O	39:BE:82:ARG:O	2.27	0.53
40:BF:34:TRP:HB3	46:BP:11:GLY:HA3	1.89	0.53
41:BG:96:ARG:HA	41:BG:99:MET:HE1	1.90	0.53
42:BH:97:ARG:HD2	42:BH:104:GLU:OE1	2.09	0.53
35:BA:598:G:H5'	46:BP:15:ARG:CB	2.38	0.53
46:BP:16:ARG:HD3	46:BP:18:ARG:N	2.20	0.53
47:BQ:81:VAL:HG23	47:BQ:82:ARG:HG2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BR:2:ARG:C	48:BR:2:ARG:HD2	2.29	0.53
48:BR:73:VAL:O	48:BR:76:VAL:HG12	2.07	0.53
48:BR:52:ILE:HD12	48:BR:79:LEU:HD21	1.90	0.53
48:BR:85:PRO:O	48:BR:87:TYR:N	2.42	0.53
50:BT:23:ARG:NH2	50:BT:120:ARG:HD3	2.23	0.53
52:BV:33:VAL:HG13	52:BV:62:LEU:H	1.72	0.53
52:BV:72:VAL:CA	52:BV:88:ARG:HH12	2.16	0.53
56:BZ:56:VAL:HA	56:BZ:70:LEU:CD2	2.38	0.53
1:CA:1372:U:H5''	9:CI:71:SER:CB	2.36	0.53
1:CA:1424:C:O2'	1:CA:1425:U:H5'	2.08	0.53
1:CA:1511:G:H2'	1:CA:1512:U:O4'	2.08	0.53
1:CA:158:G:O2'	1:CA:159:G:H5'	2.07	0.53
1:CA:15:G:H4'	5:CE:24:ARG:HH21	1.73	0.53
1:CA:328:C:H4'	1:CA:329:A:C5'	2.36	0.53
1:CA:764:C:C2'	1:CA:765:G:H8	2.19	0.53
1:CA:973:G:C4	10:CJ:55:LYS:NZ	2.68	0.53
1:CA:973:G:O4'	10:CJ:55:LYS:HG2	2.08	0.53
2:CB:72:GLY:HA3	2:CB:165:VAL:CG2	2.38	0.53
4:CD:58:LEU:O	4:CD:60:GLU:N	2.41	0.53
7:CG:22:LEU:O	7:CG:25:ALA:HB3	2.08	0.53
9:CI:7:THR:N	9:CI:83:ARG:HD2	2.24	0.53
11:CK:122:LYS:O	11:CK:126:ARG:CB	2.57	0.53
11:CK:85:ARG:HG2	11:CK:111:ASP:O	2.08	0.53
1:CA:995:C:O2	14:CN:4:LYS:HE2	2.09	0.53
19:CS:22:LEU:C	19:CS:24:ALA:H	2.12	0.53
25:CY:127:VAL:O	25:CY:128:ALA:C	2.45	0.53
25:CY:176:ALA:O	25:CY:180:GLU:HG3	2.09	0.53
25:CY:18:LEU:HD12	25:CY:18:LEU:C	2.29	0.53
27:D1:23:LYS:NZ	27:D1:23:LYS:HA	2.23	0.53
29:D3:26:LEU:HD21	29:D3:46:ASN:CB	2.38	0.53
34:D8:56:GLU:HA	34:D8:59:LYS:NZ	2.22	0.53
35:DA:1160:G:N2	52:DV:10:LYS:HE3	2.24	0.53
35:DA:1779:U:H5	35:DA:1784:A:N7	2.04	0.53
35:DA:118:A:N3	35:DA:178:G:H1'	2.23	0.53
35:DA:1856:G:H2'	35:DA:1857:G:H5'	1.89	0.53
25:CY:119:ARG:HD3	35:DA:1947:C:OP1	2.08	0.53
35:DA:2019:A:O3'	51:DU:27:LEU:HD12	2.08	0.53
35:DA:2033:A:O2'	35:DA:2034:U:P	2.66	0.53
35:DA:2065:C:H2'	35:DA:2066:C:C6	2.43	0.53
35:DA:2277:G:OP1	47:DQ:85:LYS:HB3	2.09	0.53
35:DA:235:U:H2'	35:DA:236:C:H6	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2593:U:H2'	35:DA:2594:C:C6	2.44	0.53
35:DA:493:G:C2'	35:DA:494:G:H5''	2.38	0.53
36:DB:57:A:O4'	41:DG:30:GLU:HB3	2.08	0.53
38:DD:40:THR:HG22	38:DD:41:GLY:O	2.09	0.53
39:DE:59:VAL:CG1	39:DE:63:LEU:HG	2.37	0.53
35:DA:607:U:OP1	40:DF:103:LYS:N	2.41	0.53
41:DG:59:GLU:OE1	41:DG:60:LEU:HD23	2.08	0.53
44:DN:78:TYR:CG	44:DN:79:PRO:HD3	2.43	0.53
46:DP:123:LEU:O	46:DP:125:VAL:HG12	2.07	0.53
47:DQ:35:VAL:HG22	47:DQ:101:ARG:O	2.08	0.53
48:DR:12:ARG:HD3	48:DR:16:HIS:CD2	2.43	0.53
48:DR:13:HIS:ND1	48:DR:13:HIS:C	2.62	0.53
49:DS:92:TYR:O	49:DS:93:LYS:HB2	2.09	0.53
50:DT:81:PRO:O	50:DT:82:LEU:HD12	2.08	0.53
55:DY:31:LEU:HD11	55:DY:34:LYS:H	1.72	0.53
55:DY:67:LEU:HD12	55:DY:68:HIS:N	2.23	0.53
56:DZ:87:ASP:N	56:DZ:87:ASP:OD2	2.40	0.53
1:AA:1089:G:O2'	1:AA:1090:U:H5'	2.08	0.53
1:AA:1118:C:H1'	1:AA:1179:A:C4	2.42	0.53
1:AA:1154:G:H2'	1:AA:1155:G:H8	1.73	0.53
1:AA:1347:G:C2	1:AA:1373:G:H2'	2.43	0.53
1:AA:1511:G:O2'	1:AA:1512:U:H5'	2.09	0.53
1:AA:1527:C:H2'	1:AA:1528:U:C6	2.44	0.53
1:AA:501:C:H2'	1:AA:502:G:C8	2.42	0.53
1:AA:625:G:O2'	1:AA:626:U:H5'	2.08	0.53
1:AA:967:C:H2'	1:AA:968:A:C8	2.43	0.53
1:AA:998:G:H2'	1:AA:999:C:C6	2.43	0.53
2:AB:28:PHE:O	2:AB:28:PHE:CD1	2.61	0.53
2:AB:34:ALA:O	2:AB:41:ILE:HB	2.09	0.53
2:AB:82:ARG:HG3	2:AB:92:TYR:OH	2.07	0.53
3:AC:47:LEU:HD21	3:AC:52:LEU:HD13	1.90	0.53
4:AD:131:ARG:O	4:AD:132:ARG:C	2.46	0.53
4:AD:92:VAL:O	4:AD:95:GLY:N	2.41	0.53
5:AE:12:LEU:O	5:AE:12:LEU:HD22	2.08	0.53
5:AE:147:ASP:CA	5:AE:150:ARG:HB3	2.37	0.53
6:AF:29:ALA:HB1	6:AF:79:LEU:CD2	2.39	0.53
7:AG:105:VAL:HG12	7:AG:109:ASN:HD21	1.72	0.53
8:AH:119:LEU:N	8:AH:119:LEU:CD2	2.63	0.53
11:AK:109:VAL:HG13	18:AR:85:LEU:O	2.09	0.53
13:AM:108:ARG:CZ	13:AM:114:ARG:HG2	2.39	0.53
14:AN:41:ARG:HG3	14:AN:42:ILE:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AN:6:LEU:O	14:AN:8:GLU:N	2.42	0.53
16:AP:23:ASP:HB3	16:AP:26:ARG:HG3	1.89	0.53
16:AP:72:ARG:O	16:AP:74:LEU:N	2.41	0.53
17:AQ:48:GLU:O	17:AQ:49:GLU:C	2.47	0.53
17:AQ:92:ARG:O	17:AQ:95:TYR:N	2.35	0.53
19:AS:64:GLU:HG3	19:AS:65:ASN:N	2.22	0.53
1:AA:1243:C:OP2	21:AU:10:ARG:CZ	2.56	0.53
23:AW:12:G:H1	23:AW:24:C:H42	1.57	0.53
1:AA:693:G:H22	23:AW:38:A:H2	1.55	0.53
25:AY:70:SER:OG	25:AY:76:LEU:HB2	2.08	0.53
27:B1:11:ARG:O	27:B1:13:ILE:N	2.41	0.53
27:B1:18:ILE:HG12	27:B1:43:TYR:CD1	2.43	0.53
28:B2:32:LEU:HA	28:B2:37:PHE:CD2	2.43	0.53
35:BA:1159:U:C2'	35:BA:1160:G:H5'	2.37	0.53
35:BA:1591:G:O2'	35:BA:1592:C:H5'	2.08	0.53
35:BA:204:A:H5'	35:BA:206:U:O4'	2.07	0.53
35:BA:2114:A:H2'	35:BA:2114:A:N3	2.24	0.53
35:BA:2620:C:OP1	39:BE:152:LYS:O	2.27	0.53
35:BA:2692:C:H2'	35:BA:2693:A:H8	1.73	0.53
35:BA:2773:C:H2'	35:BA:2774:C:H6	1.73	0.53
35:BA:468:G:H2'	35:BA:469:G:O4'	2.07	0.53
35:BA:477:A:H2'	35:BA:478:A:C8	2.43	0.53
35:BA:834:C:H2'	35:BA:835:A:H8	1.72	0.53
35:BA:847:U:C2'	35:BA:848:G:H5''	2.30	0.53
38:BD:227:ASN:O	38:BD:229:VAL:N	2.42	0.53
39:BE:36:ARG:NH2	39:BE:88:GLY:H	2.05	0.53
41:BG:114:ILE:HB	41:BG:117:PHE:HB2	1.89	0.53
42:BH:19:VAL:HG11	42:BH:44:VAL:HG22	1.91	0.53
44:BN:70:LYS:HB3	44:BN:87:LEU:HB2	1.91	0.53
44:BN:7:LYS:O	44:BN:9:VAL:N	2.41	0.53
44:BN:97:ARG:O	44:BN:101:HIS:N	2.36	0.53
46:BP:101:VAL:HG13	46:BP:102:ARG:N	2.21	0.53
46:BP:124:LYS:HA	46:BP:143:GLY:CA	2.38	0.53
47:BQ:42:ILE:HD13	47:BQ:97:VAL:HG21	1.89	0.53
48:BR:12:ARG:HD3	48:BR:16:HIS:CD2	2.43	0.53
50:BT:81:PRO:C	50:BT:82:LEU:HD12	2.28	0.53
47:BQ:141:GLN:C	56:BZ:53:ILE:HB	2.29	0.53
1:CA:261:U:H3'	20:CT:79:ARG:HH12	1.74	0.53
1:CA:357:G:OP1	1:CA:366:C:O2'	2.25	0.53
1:CA:451:A:N6	1:CA:480:U:H2'	2.23	0.53
1:CA:882:C:H2'	1:CA:883:C:H6	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:171:ALA:O	2:CB:174:VAL:N	2.41	0.53
2:CB:73:THR:HB	2:CB:94:ASN:O	2.08	0.53
4:CD:176:LEU:HD21	4:CD:178:VAL:HG22	1.90	0.53
8:CH:83:ILE:HD12	8:CH:136:GLU:O	2.08	0.53
10:CJ:16:LEU:HD22	10:CJ:19:SER:OG	2.09	0.53
11:CK:99:GLN:NE2	11:CK:105:VAL:HG11	2.23	0.53
13:CM:35:GLU:HG3	13:CM:36:LYS:N	2.23	0.53
15:CO:39:LEU:O	15:CO:39:LEU:HD13	2.09	0.53
18:CR:59:SER:OG	18:CR:60:ALA:N	2.40	0.53
25:CY:134:ARG:O	25:CY:135:GLU:C	2.46	0.53
25:CY:166:ASP:O	25:CY:167:GLU:C	2.46	0.53
27:D1:16:ASN:N	27:D1:16:ASN:HD22	1.98	0.53
28:D2:25:VAL:O	28:D2:26:ARG:C	2.47	0.53
28:D2:40:SER:HB3	28:D2:41:ILE:HD12	1.88	0.53
35:DA:1174:A:H5''	35:DA:1175:U:H5''	1.90	0.53
35:DA:1245:G:H3'	46:DP:16:ARG:NH2	2.22	0.53
35:DA:1386:C:O2'	35:DA:1387:C:H5'	2.08	0.53
35:DA:1414:G:H2'	35:DA:1415:U:C5	2.43	0.53
35:DA:1902:C:H5'	38:DD:246:PRO:HD3	1.91	0.53
35:DA:2022:U:O2'	35:DA:2617:C:H5'	2.08	0.53
35:DA:2059:A:H5'	35:DA:2060:A:OP2	2.08	0.53
26:D0:43:THR:N	35:DA:2331:G:H4'	2.23	0.53
35:DA:2752:C:H5'	35:DA:2753:A:OP2	2.07	0.53
35:DA:62:C:H2'	35:DA:63:U:H5'	1.89	0.53
38:DD:118:VAL:CG2	38:DD:119:ALA:N	2.71	0.53
38:DD:130:ALA:CB	38:DD:192:THR:HA	2.38	0.53
38:DD:5:LYS:N	38:DD:5:LYS:HD2	2.23	0.53
39:DE:179:GLU:OE1	39:DE:179:GLU:HA	2.08	0.53
35:DA:2632:A:H1'	39:DE:61:ARG:NH1	2.23	0.53
41:DG:108:ASN:O	41:DG:112:PRO:HB2	2.08	0.53
43:DI:120:ILE:O	43:DI:121:LYS:HB3	2.08	0.53
43:DI:41:GLU:HA	43:DI:44:LEU:HB3	1.90	0.53
45:DO:12:ASP:OD2	45:DO:12:ASP:N	2.39	0.53
49:DS:95:HIS:O	49:DS:97:ARG:N	2.41	0.53
51:DU:92:ARG:CZ	52:DV:11:GLN:HG2	2.39	0.53
52:DV:34:GLU:C	52:DV:62:LEU:HD12	2.27	0.53
53:DW:64:MET:HE2	53:DW:109:GLU:HG3	1.90	0.53
53:DW:86:LEU:HD12	53:DW:87:PRO:HD2	1.90	0.53
54:DX:29:TRP:HA	54:DX:29:TRP:CE3	2.43	0.53
55:DY:98:VAL:HG12	55:DY:98:VAL:O	2.09	0.53
56:DZ:131:ARG:C	56:DZ:133:ILE:HD12	2.28	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1057:G:H5'	3:AC:154:SER:OG	2.07	0.53
1:AA:1244:C:H2'	1:AA:1245:A:H8	1.71	0.53
1:AA:486:U:O2'	1:AA:487:A:H5'	2.08	0.53
4:AD:129:ASN:HB2	4:AD:131:ARG:NH2	2.24	0.53
4:AD:8:VAL:O	4:AD:9:CYS:C	2.46	0.53
8:AH:83:ILE:HD12	8:AH:136:GLU:O	2.07	0.53
11:AK:49:GLY:O	11:AK:50:TYR:HD2	1.91	0.53
16:AP:67:THR:HB	16:AP:70:ALA:H	1.74	0.53
17:AQ:10:VAL:HG21	17:AQ:55:ASP:HB2	1.90	0.53
19:AS:9:VAL:O	19:AS:11:VAL:N	2.41	0.53
19:AS:16:LEU:N	19:AS:16:LEU:HD12	2.24	0.53
20:AT:38:LYS:O	20:AT:41:ILE:N	2.41	0.53
20:AT:63:ILE:O	20:AT:66:ALA:N	2.42	0.53
23:AW:53:G:N2	23:AW:64:G:N1	2.57	0.53
27:B1:24:ALA:HA	27:B1:36:GLY:HA2	1.89	0.53
28:B2:49:LYS:O	28:B2:53:LEU:O	2.25	0.53
31:B5:11:THR:O	31:B5:12:SER:C	2.46	0.53
33:B7:34:ARG:C	33:B7:36:GLN:N	2.62	0.53
34:B8:62:LEU:HD13	35:BA:242:G:C5'	2.20	0.53
35:BA:1380:G:C2	35:BA:1381:G:C8	2.97	0.53
35:BA:221:A:O2'	35:BA:222:A:OP2	2.23	0.53
35:BA:2442:C:O2'	35:BA:2443:C:H5'	2.08	0.53
35:BA:2642:G:N2	35:BA:2773:C:C2	2.77	0.53
35:BA:301:G:C4	35:BA:302:C:C5	2.96	0.53
35:BA:590:A:H2'	35:BA:591:C:H6	1.73	0.53
35:BA:668:G:H2'	35:BA:670:A:H62	1.74	0.53
35:BA:587:C:C6	35:BA:671:C:H1'	2.44	0.53
35:BA:685:A:N1	35:BA:787:U:H1'	2.24	0.53
37:BC:23:ASP:C	37:BC:25:ALA:H	2.12	0.53
37:BC:82:LYS:O	37:BC:86:ALA:HB3	2.08	0.53
39:BE:117:MET:CE	39:BE:124:GLY:HA3	2.38	0.53
39:BE:69:LYS:N	39:BE:69:LYS:HE2	2.23	0.53
40:BF:110:LEU:HA	40:BF:183:VAL:CG1	2.36	0.53
42:BH:27:LYS:HG2	42:BH:32:GLU:OE1	2.08	0.53
42:BH:67:LEU:O	42:BH:71:LEU:HD22	2.07	0.53
42:BH:87:LEU:HD23	42:BH:164:TYR:HA	1.91	0.53
43:BI:76:THR:HB	43:BI:139:GLN:O	2.08	0.53
35:BA:1190:G:O3'	46:BP:35:HIS:HB3	2.08	0.53
47:BQ:76:LYS:H	47:BQ:88:GLY:HA3	1.72	0.53
47:BQ:8:LYS:CG	47:BQ:9:TYR:N	2.64	0.53
48:BR:16:HIS:O	48:BR:17:ARG:C	2.47	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:BU:90:VAL:HG13	52:BV:39:LEU:CB	2.34	0.53
52:BV:4:ILE:HA	52:BV:12:TYR:O	2.08	0.53
56:BZ:33:LEU:HD11	56:BZ:35:ARG:CG	2.39	0.53
1:CA:1006:C:O2'	1:CA:1007:C:H5'	2.08	0.53
1:CA:102:G:H2'	1:CA:103:C:C6	2.43	0.53
1:CA:1043:C:H2'	1:CA:1044:A:C8	2.43	0.53
1:CA:1072:G:H2'	1:CA:1073:U:H6	1.70	0.53
1:CA:1242:C:O5'	1:CA:1242:C:H6	1.92	0.53
1:CA:397:A:H5'	1:CA:398:C:OP1	2.08	0.53
1:CA:582:U:H2'	1:CA:583:A:H8	1.73	0.53
2:CB:114:ARG:HA	2:CB:117:GLU:OE1	2.09	0.53
2:CB:208:ILE:O	2:CB:208:ILE:HG22	2.07	0.53
2:CB:80:ILE:HD13	2:CB:212:GLN:HA	1.89	0.53
2:CB:61:LEU:O	2:CB:64:ARG:HG2	2.09	0.53
1:CA:1112:C:C4	3:CC:178:LEU:HD23	2.43	0.53
15:CO:12:ILE:C	15:CO:14:GLU:H	2.11	0.53
19:CS:43:GLU:O	19:CS:45:VAL:N	2.39	0.53
19:CS:42:PRO:C	19:CS:44:MET:H	2.10	0.53
20:CT:38:LYS:O	20:CT:41:ILE:N	2.41	0.53
25:CY:162:GLN:O	25:CY:165:THR:HB	2.07	0.53
27:D1:43:TYR:OH	35:DA:1365:A:H5'	2.09	0.53
27:D1:89:GLU:HG2	27:D1:90:ILE:HD13	1.91	0.53
28:D2:23:LYS:HA	54:DX:5:TYR:CE1	2.38	0.53
29:D3:21:ALA:O	29:D3:24:LYS:N	2.39	0.53
29:D3:3:ARG:HG2	29:D3:38:GLU:OE2	2.08	0.53
29:D3:41:PRO:HD3	29:D3:44:ARG:NH1	2.23	0.53
31:D5:11:THR:HB	35:DA:1263:U:O3'	2.09	0.53
32:D6:30:THR:HG21	35:DA:2286:A:OP1	2.08	0.53
34:D8:31:HIS:O	34:D8:32:LEU:C	2.47	0.53
34:D8:39:LYS:HD3	34:D8:39:LYS:C	2.28	0.53
35:DA:1289:C:H2'	35:DA:1290:C:H6	1.74	0.53
35:DA:1469:A:H2'	35:DA:1470:G:O4'	2.09	0.53
35:DA:1475:G:N3	35:DA:1475:G:H2'	2.23	0.53
35:DA:2039:C:H2'	35:DA:2040:C:H6	1.72	0.53
35:DA:2248:C:O2'	35:DA:2249:U:H5'	2.08	0.53
35:DA:1786:A:N1	35:DA:2606:C:H1'	2.24	0.53
35:DA:2748:A:C2	42:DH:63:SER:HB3	2.44	0.53
35:DA:2807:G:C3'	35:DA:2808:U:H5''	2.38	0.53
35:DA:2839:G:H2'	35:DA:2840:C:C6	2.43	0.53
35:DA:590:A:H2'	35:DA:591:C:H6	1.74	0.53
35:DA:848:G:H2'	35:DA:849:A:C8	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:921:G:H2'	35:DA:922:U:H6	1.72	0.53
36:DB:65:C:H2'	36:DB:66:A:H5'	1.90	0.53
38:DD:130:ALA:HB2	38:DD:192:THR:HA	1.91	0.53
38:DD:9:TYR:CD2	38:DD:10:THR:HG22	2.43	0.53
40:DF:53:THR:O	40:DF:54:ARG:C	2.46	0.53
46:DP:50:ARG:HD2	46:DP:51:PHE:CG	2.43	0.53
46:DP:58:THR:O	46:DP:61:ARG:NE	2.42	0.53
35:DA:2406:U:C4	46:DP:72:PRO:HD2	2.43	0.53
35:DA:2406:U:N3	46:DP:72:PRO:HD2	2.22	0.53
49:DS:42:ASP:O	49:DS:44:LYS:N	2.42	0.53
49:DS:65:VAL:CG1	49:DS:69:VAL:HB	2.39	0.53
51:DU:31:SER:C	51:DU:33:ARG:N	2.62	0.53
52:DV:15:GLU:HB3	52:DV:16:PRO:CD	2.37	0.53
54:DX:31:HIS:ND1	54:DX:32:PRO:HD2	2.23	0.53
1:AA:1287:A:C6	1:AA:1288:A:N6	2.76	0.53
1:AA:1239:A:N6	1:AA:1299:A:H62	2.05	0.53
1:AA:1347:G:O2'	1:AA:1348:U:OP2	2.27	0.53
1:AA:741:G:H5'	15:AO:39:LEU:CD2	2.39	0.53
2:AB:75:LYS:HA	2:AB:75:LYS:HE3	1.91	0.53
3:AC:16:ARG:NH1	3:AC:16:ARG:CB	2.69	0.53
3:AC:83:ARG:O	3:AC:86:VAL:HG22	2.08	0.53
5:AE:92:LYS:HB2	5:AE:119:LEU:HB2	1.91	0.53
9:AI:26:VAL:HG13	9:AI:63:ILE:HD11	1.90	0.53
11:AK:18:ARG:HG2	11:AK:33:THR:OG1	2.09	0.53
12:AL:24:VAL:O	12:AL:24:VAL:HG12	2.08	0.53
13:AM:67:GLU:O	13:AM:69:GLU:N	2.42	0.53
1:AA:656:C:H4'	15:AO:62:GLN:NE2	2.23	0.53
16:AP:49:LEU:HD11	16:AP:51:VAL:CG2	2.35	0.53
23:AW:60:A:O2'	23:AW:61:U:H5'	2.08	0.53
25:AY:66:LEU:O	25:AY:101:ILE:HD12	2.08	0.53
27:B1:67:ILE:N	27:B1:68:PRO:HD2	2.23	0.53
27:B1:75:GLU:O	27:B1:76:ARG:NH2	2.41	0.53
30:B4:1:MET:N	36:BB:43:C:H5'	2.23	0.53
31:B5:31:VAL:HB	31:B5:32:PRO:CD	2.32	0.53
35:BA:1153:C:N4	35:BA:1154:G:C2	2.76	0.53
35:BA:1452:A:O2'	35:BA:1453:U:H2'	2.07	0.53
35:BA:1467:C:OP2	35:BA:1547:C:H5	1.91	0.53
35:BA:1478:G:HO2'	35:BA:1558:A:H2	1.55	0.53
35:BA:210:C:O2'	35:BA:211:A:H5'	2.08	0.53
35:BA:2649:U:O2'	35:BA:2650:U:H5'	2.09	0.53
38:BD:247:ALA:CA	38:BD:254:THR:HG22	2.35	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1812:A:H1'	38:BD:46:GLN:HE22	1.74	0.53
40:BF:183:VAL:O	40:BF:187:VAL:HG23	2.08	0.53
40:BF:57:VAL:CG1	40:BF:58:ALA:N	2.71	0.53
42:BH:118:PRO:HG2	42:BH:121:ILE:HD12	1.90	0.53
44:BN:42:TRP:CE3	44:BN:42:TRP:HA	2.44	0.53
46:BP:46:LYS:CB	46:BP:52:GLU:HG2	2.38	0.53
48:BR:87:TYR:OH	48:BR:116:LEU:HD22	2.09	0.53
49:BS:12:PHE:CG	49:BS:12:PHE:O	2.62	0.53
50:BT:96:ARG:HG2	50:BT:96:ARG:NH1	2.20	0.53
52:BV:22:VAL:HG21	52:BV:96:ILE:HB	1.91	0.53
55:BY:83:THR:O	55:BY:84:ARG:HG3	2.08	0.53
56:BZ:7:ALA:O	56:BZ:8:TYR:O	2.26	0.53
1:CA:1221:G:OP1	19:CS:36:ARG:HD3	2.08	0.53
1:CA:511:C:C2	1:CA:512:U:C5	2.96	0.53
1:CA:738:C:H5''	6:CF:69:GLU:CB	2.39	0.53
1:CA:908:A:C4	1:CA:909:A:N7	2.76	0.53
2:CB:187:LEU:HA	2:CB:201:ILE:O	2.08	0.53
2:CB:87:ARG:O	2:CB:88:ALA:HB2	2.08	0.53
4:CD:8:VAL:O	4:CD:9:CYS:C	2.47	0.53
7:CG:103:TRP:O	7:CG:104:LEU:C	2.46	0.53
7:CG:120:ILE:C	7:CG:124:LEU:HD12	2.29	0.53
8:CH:83:ILE:CB	8:CH:137:VAL:HG13	2.36	0.53
9:CI:11:LYS:O	9:CI:13:ALA:N	2.42	0.53
9:CI:16:ARG:O	9:CI:63:ILE:CG2	2.56	0.53
9:CI:65:VAL:HG22	9:CI:66:ARG:H	1.73	0.53
11:CK:21:ILE:HD13	11:CK:84:VAL:CG1	2.39	0.53
11:CK:18:ARG:HG2	11:CK:33:THR:OG1	2.08	0.53
11:CK:62:GLN:C	11:CK:64:ALA:N	2.61	0.53
13:CM:13:LYS:HB2	13:CM:18:ALA:HB2	1.91	0.53
16:CP:67:THR:HB	16:CP:70:ALA:H	1.73	0.53
25:CY:165:THR:O	25:CY:166:ASP:C	2.47	0.53
32:D6:9:LEU:HD23	32:D6:10:LEU:O	2.08	0.53
35:DA:1019:U:H3	35:DA:1142(A):A:N6	1.98	0.53
35:DA:1354:A:C8	35:DA:1355:G:C8	2.97	0.53
35:DA:1453:U:H5'	48:DR:63:ARG:HE	1.72	0.53
35:DA:1678:G:N2	35:DA:1989:G:H22	2.05	0.53
35:DA:2181:G:H2'	35:DA:2182:G:C8	2.44	0.53
35:DA:2574:G:H2'	35:DA:2575:C:C6	2.42	0.53
35:DA:2683:C:H5''	50:DT:53:ARG:NH2	2.22	0.53
35:DA:2781:A:H5'	35:DA:2782:G:H5'	1.87	0.53
35:DA:304:G:H1	35:DA:313:C:H42	1.55	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:589:C:H42	35:DA:668:G:H1	1.56	0.53
35:DA:815:C:OP2	52:DV:84:LYS:HE3	2.09	0.53
37:DC:82:LYS:O	37:DC:86:ALA:HB3	2.08	0.53
35:DA:782:A:H2	38:DD:226:MET:HG2	1.66	0.53
35:DA:2619:C:H5''	39:DE:152:LYS:HA	1.90	0.53
41:DG:43:LEU:HD21	41:DG:88:ILE:HG22	1.91	0.53
43:DI:17:GLN:HG2	43:DI:18:VAL:N	2.21	0.53
44:DN:26:LEU:HG	44:DN:30:ILE:CD1	2.39	0.53
46:DP:127:ALA:HB3	46:DP:130:PHE:HE2	1.64	0.53
47:DQ:17:LEU:HD21	47:DQ:41:TRP:HE1	1.73	0.53
47:DQ:28:ALA:HB1	47:DQ:29:PHE:CD1	2.44	0.53
49:DS:26:LEU:HD13	49:DS:87:PHE:HD1	1.74	0.53
50:DT:101:PHE:CD2	50:DT:102:ILE:N	2.71	0.53
50:DT:38:ASN:ND2	50:DT:40:THR:OG1	2.42	0.53
51:DU:64:ARG:NH2	51:DU:64:ARG:HA	2.23	0.53
56:DZ:53:ILE:HG22	56:DZ:71:VAL:CB	2.38	0.53
56:DZ:76:LEU:HD22	56:DZ:76:LEU:N	2.23	0.53
1:AA:1132:C:N4	1:AA:1133:G:C6	2.77	0.53
1:AA:1419:G:N2	1:AA:1482:G:C4	2.76	0.53
1:AA:176:C:H2'	1:AA:177:C:C6	2.43	0.53
1:AA:865:A:H2'	1:AA:866:C:C6	2.43	0.53
1:AA:959:A:H2'	1:AA:960:U:C4'	2.39	0.53
2:AB:208:ILE:HG22	2:AB:208:ILE:O	2.09	0.53
3:AC:3:ASN:O	3:AC:4:LYS:C	2.47	0.53
5:AE:8:GLU:CA	5:AE:34:VAL:HG22	2.35	0.53
6:AF:37:VAL:HA	6:AF:65:VAL:CG1	2.38	0.53
8:AH:34:GLU:O	8:AH:38:ILE:HG13	2.09	0.53
8:AH:83:ILE:HG23	8:AH:83:ILE:O	2.08	0.53
10:AJ:34:VAL:HG12	10:AJ:35:SER:H	1.72	0.53
10:AJ:57:LYS:HD2	10:AJ:60:ARG:NH2	2.23	0.53
11:AK:103:LEU:N	11:AK:103:LEU:CD2	2.65	0.53
1:AA:280:C:O2	17:AQ:38:ARG:HG3	2.09	0.53
19:AS:33:THR:HG21	19:AS:51:VAL:HG22	1.90	0.53
20:AT:81:LYS:O	20:AT:85:MET:HG2	2.09	0.53
25:AY:156:ARG:HH21	26:B0:6:ALA:HB2	1.74	0.53
25:AY:29:ARG:HE	25:AY:32:ARG:HH22	1.54	0.53
27:B1:87:PRO:HB2	27:B1:91:LYS:HE3	1.91	0.53
28:B2:27:GLU:C	28:B2:29:LYS:H	2.11	0.53
28:B2:43:GLN:O	28:B2:46:GLN:OE1	2.26	0.53
33:B7:31:LEU:CD2	33:B7:42:LEU:HB3	2.38	0.53
35:BA:1021:A:C8	35:BA:1021:A:H3'	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1567:A:H5'	38:BD:58:HIS:CD2	2.44	0.53
35:BA:1830:C:N4	35:BA:1975:G:H1	2.07	0.53
35:BA:1917:U:H2'	35:BA:1918:A:C8	2.44	0.53
35:BA:2009:G:C6	35:BA:2010:G:N7	2.76	0.53
35:BA:2704:C:C2'	35:BA:2705:A:H8	2.17	0.53
35:BA:524:U:O2	35:BA:524:U:C2'	2.53	0.53
35:BA:570:G:H2'	35:BA:2030:A:C5	2.43	0.53
35:BA:796:C:H2'	35:BA:797:C:H6	1.72	0.53
35:BA:941:A:H4'	46:BP:35:HIS:CE1	2.44	0.53
36:BB:31:C:H42	36:BB:51:G:H1	1.55	0.53
38:BD:262:ARG:O	38:BD:264:LYS:N	2.41	0.53
39:BE:3:GLY:O	39:BE:4:ILE:HG22	2.08	0.53
42:BH:85:LYS:NZ	42:BH:145:ALA:HA	2.24	0.53
45:BO:37:ASP:O	45:BO:61:VAL:HA	2.09	0.53
35:BA:1667:G:OP1	45:BO:7:TYR:HB2	2.08	0.53
45:BO:88:ASN:OD1	45:BO:92:GLU:N	2.36	0.53
46:BP:17:LYS:C	46:BP:19:VAL:N	2.61	0.53
35:BA:1030:G:OP2	47:BQ:128:LYS:HE2	2.09	0.53
47:BQ:28:ALA:HB3	47:BQ:105:GLU:CD	2.29	0.53
48:BR:49:ASP:O	48:BR:51:LEU:N	2.41	0.53
50:BT:36:GLU:HB3	50:BT:38:ASN:OD1	2.08	0.53
52:BV:27:ALA:O	52:BV:29:PRO:N	2.42	0.53
35:BA:519:U:H4'	53:BW:25:ARG:HH22	1.74	0.53
55:BY:37:VAL:HG13	55:BY:69:ALA:CB	2.39	0.53
56:BZ:44:PHE:CD1	56:BZ:48:PHE:HB2	2.44	0.53
1:CA:1023:G:H2'	1:CA:1024:G:H5'	1.91	0.53
1:CA:1132:C:N4	1:CA:1133:G:C6	2.76	0.53
1:CA:236:G:H2'	1:CA:237:C:H6	1.72	0.53
1:CA:116:A:H61	1:CA:313:A:H1'	1.74	0.53
1:CA:645:C:O2'	1:CA:646:U:H5'	2.09	0.53
1:CA:856:C:H2'	1:CA:857:C:H6	1.74	0.53
2:CB:82:ARG:HG3	2:CB:92:TYR:OH	2.09	0.53
3:CC:52:LEU:O	3:CC:53:ALA:HB2	2.09	0.53
4:CD:93:PHE:CE1	4:CD:97:LEU:HD12	2.43	0.53
4:CD:96:LEU:C	4:CD:98:GLU:N	2.62	0.53
5:CE:15:ARG:HG2	5:CE:26:PHE:HD2	1.74	0.53
7:CG:58:PRO:HA	7:CG:61:VAL:HG23	1.91	0.53
12:CL:102:ARG:CG	12:CL:102:ARG:HH11	2.18	0.53
13:CM:108:ARG:CZ	13:CM:114:ARG:HG2	2.38	0.53
13:CM:74:VAL:O	13:CM:78:ILE:HG13	2.08	0.53
23:CW:25:U:H2'	23:CW:26:C:C6	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:CX:12:G:H3'	24:CX:13:A:H5''	1.90	0.53
35:DA:1378:A:C4'	35:DA:1379:A:OP1	2.55	0.53
35:DA:149:A:H2'	35:DA:150:C:C6	2.42	0.53
35:DA:1553:A:H2'	35:DA:1554:A:H5''	1.89	0.53
35:DA:16:G:H2'	35:DA:17:G:H8	1.73	0.53
35:DA:2547:U:O2'	35:DA:2548:G:H5'	2.09	0.53
35:DA:2545:G:N3	35:DA:2565:A:H2	2.07	0.53
35:DA:88:G:H2'	35:DA:88:G:N3	2.23	0.53
35:DA:967:C:C2'	35:DA:968:G:H5'	2.39	0.53
38:DD:35:LYS:HG2	38:DD:64:ILE:HG22	1.89	0.53
38:DD:65:ILE:CD1	38:DD:65:ILE:C	2.77	0.53
39:DE:1:MET:HB3	39:DE:200:GLU:OE1	2.09	0.53
39:DE:36:ARG:NH2	39:DE:88:GLY:H	2.05	0.53
40:DF:3:GLU:HG3	40:DF:19:GLU:HB2	1.91	0.53
41:DG:63:ILE:C	41:DG:63:ILE:HD12	2.29	0.53
42:DH:27:LYS:HG2	42:DH:32:GLU:OE1	2.09	0.53
44:DN:7:LYS:O	44:DN:9:VAL:N	2.41	0.53
45:DO:37:ASP:O	45:DO:39:ILE:HG22	2.08	0.53
49:DS:57:LYS:O	49:DS:58:LEU:HB2	2.09	0.53
52:DV:78:LYS:HD3	52:DV:78:LYS:C	2.29	0.53
55:DY:8:LYS:H	55:DY:8:LYS:CD	2.09	0.53
56:DZ:103:ARG:O	56:DZ:139:VAL:HB	2.08	0.53
1:AA:1104:G:H2'	1:AA:1105:A:C8	2.44	0.53
1:AA:119:A:O2'	1:AA:120:A:OP2	2.25	0.53
1:AA:1305:G:H5''	21:AU:5:ASP:N	2.24	0.53
1:AA:1526:G:H2'	1:AA:1527:C:C6	2.44	0.53
1:AA:1526:G:H2'	1:AA:1527:C:H6	1.74	0.53
1:AA:405:U:H5''	1:AA:406:G:O4'	2.09	0.53
1:AA:727:G:N1	1:AA:731:G:C6	2.77	0.53
3:AC:70:VAL:O	3:AC:105:GLU:HA	2.09	0.53
4:AD:119:GLN:HE21	4:AD:123:HIS:CD2	2.27	0.53
5:AE:130:ASN:O	5:AE:131:ILE:C	2.47	0.53
6:AF:69:GLU:CD	6:AF:69:GLU:H	2.11	0.53
1:AA:939:G:H5''	7:AG:102:ARG:NH1	2.23	0.53
7:AG:58:PRO:HA	7:AG:61:VAL:HG23	1.89	0.53
8:AH:83:ILE:HB	8:AH:137:VAL:CG1	2.37	0.53
9:AI:24:GLY:HA2	9:AI:59:PHE:O	2.08	0.53
1:AA:963:G:N2	10:AJ:55:LYS:NZ	2.57	0.53
14:AN:51:GLY:O	14:AN:53:LEU:N	2.39	0.53
15:AO:11:VAL:HG13	15:AO:15:PHE:HE1	1.74	0.53
19:AS:36:ARG:CZ	19:AS:72:GLY:HA2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AT:54:LYS:HA	20:AT:57:ARG:NH1	2.24	0.53
23:AW:16:C:H6	23:AW:60:A:C2	2.27	0.53
25:AY:3:LEU:N	25:AY:3:LEU:HD12	2.16	0.53
27:B1:56:GLN:O	27:B1:57:GLU:CG	2.53	0.53
33:B7:30:VAL:O	33:B7:34:ARG:N	2.38	0.53
35:BA:1257:C:C2	35:BA:1258:C:C5	2.96	0.53
35:BA:1993:U:O2'	35:BA:1994:C:H5'	2.08	0.53
35:BA:2012:G:OP1	53:BW:98:LYS:HA	2.08	0.53
35:BA:2079:U:H3	35:BA:2241:A:N6	2.05	0.53
35:BA:2694:G:O2'	35:BA:2695:C:H5'	2.09	0.53
35:BA:2821:A:OP2	48:BR:2:ARG:NH1	2.42	0.53
35:BA:304:G:H1	35:BA:313:C:H42	1.54	0.53
35:BA:363(A):A:C2	35:BA:363(B):G:C8	2.97	0.53
35:BA:513:A:H1'	51:BU:11:ARG:NH1	2.23	0.53
35:BA:674:G:O2'	40:BF:74:ARG:HB2	2.09	0.53
37:BC:65:PRO:HG2	37:BC:189:ILE:HA	1.90	0.53
38:BD:182:LEU:O	38:BD:271:ILE:HD12	2.09	0.53
39:BE:44:TYR:O	39:BE:45:THR:CB	2.55	0.53
39:BE:59:VAL:O	39:BE:62:PRO:HD2	2.09	0.53
39:BE:6:GLY:HA2	39:BE:51:PHE:HE2	1.71	0.53
41:BG:85:GLY:O	41:BG:86:MET:CB	2.57	0.53
46:BP:5:ASP:OD2	46:BP:6:LEU:HD22	2.09	0.53
50:BT:32:TYR:O	50:BT:33:LYS:CB	2.57	0.53
51:BU:3:ARG:O	51:BU:3:ARG:HG3	2.08	0.53
54:BX:32:PRO:HD3	54:BX:72:LYS:HZ2	1.71	0.53
54:BX:9:LEU:HG	54:BX:29:TRP:O	2.09	0.53
1:CA:1046:A:H3'	1:CA:1047:G:C8	2.42	0.53
1:CA:509:A:H4'	1:CA:510:A:OP1	2.08	0.53
1:CA:511:C:HO2'	1:CA:512:U:H6	1.54	0.53
4:CD:8:VAL:O	4:CD:10:ARG:HB3	2.08	0.53
4:CD:150:GLU:HA	4:CD:153:ARG:CD	2.38	0.53
6:CF:26:ILE:O	6:CF:29:ALA:HB3	2.09	0.53
7:CG:121:ALA:CA	7:CG:124:LEU:HD12	2.38	0.53
7:CG:69:VAL:HG22	7:CG:135:VAL:HG22	1.90	0.53
7:CG:24:THR:O	7:CG:25:ALA:C	2.46	0.53
10:CJ:45:ARG:HG3	10:CJ:45:ARG:HH11	1.74	0.53
10:CJ:62:HIS:H	10:CJ:62:HIS:HD2	1.53	0.53
16:CP:1:MET:SD	16:CP:1:MET:N	2.77	0.53
18:CR:59:SER:N	18:CR:62:GLU:OE1	2.42	0.53
18:CR:87:ARG:HG2	18:CR:88:LYS:N	2.23	0.53
19:CS:36:ARG:NH1	19:CS:75:ALA:HB3	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1320:C:H5'	19:CS:70:LYS:CE	2.39	0.53
20:CT:14:LYS:HE3	20:CT:18:GLN:NE2	2.23	0.53
23:CW:43:G:H2'	23:CW:44:A:O4'	2.08	0.53
23:CW:76:C:H5'	27:D1:27:GLU:OE1	2.08	0.53
27:D1:56:GLN:O	27:D1:57:GLU:CB	2.56	0.53
27:D1:72:GLU:O	27:D1:76:ARG:CZ	2.57	0.53
35:DA:1038:C:H42	35:DA:1117:G:H1	1.57	0.53
35:DA:1186:G:H2'	35:DA:1187:G:H5'	1.91	0.53
35:DA:1374:G:H2'	35:DA:1375:C:C6	2.43	0.53
35:DA:1684:C:H42	35:DA:1704:G:H1	1.56	0.53
35:DA:1714:G:H2'	35:DA:1717:G:H8	1.72	0.53
35:DA:1826:G:H4'	38:DD:242:ARG:NH2	2.24	0.53
35:DA:1918:A:O2'	35:DA:1920:C:N4	2.42	0.53
35:DA:2090:G:C6	35:DA:2091:U:C4	2.97	0.53
35:DA:2870:C:O2'	35:DA:2871:C:H5'	2.07	0.53
35:DA:668:G:H2'	35:DA:670:A:H62	1.73	0.53
35:DA:840:C:O2'	35:DA:841:A:H5'	2.08	0.53
38:DD:65:ILE:HD11	38:DD:67:PHE:CE1	2.43	0.53
35:DA:1655:A:H4'	39:DE:115:GLY:H	1.74	0.53
39:DE:59:VAL:HG13	39:DE:60:ASN:N	2.24	0.53
40:DF:151:SER:C	40:DF:152:GLU:HG3	2.28	0.53
40:DF:169:ASN:O	40:DF:169:ASN:ND2	2.42	0.53
41:DG:48:GLU:CG	41:DG:49:ASP:H	2.22	0.53
41:DG:57:ALA:O	41:DG:68:PRO:HG3	2.09	0.53
42:DH:137:ASP:O	42:DH:138:LYS:CB	2.57	0.53
42:DH:92:ILE:HG22	42:DH:93:GLY:H	1.74	0.53
43:DI:10:GLU:CD	43:DI:11:ASN:N	2.62	0.53
46:DP:70:GLN:HA	46:DP:70:GLN:OE1	2.08	0.53
47:DQ:101:ARG:HG2	47:DQ:102:VAL:N	2.23	0.53
47:DQ:73:PRO:HA	47:DQ:93:TYR:CD2	2.43	0.53
50:DT:43:GLN:HG2	50:DT:44:ASP:O	2.09	0.53
35:DA:2875:C:O2'	50:DT:5:ALA:HB3	2.09	0.53
53:DW:74:ALA:O	53:DW:75:TYR:CB	2.57	0.53
54:DX:62:LYS:CB	54:DX:69:TYR:H	2.22	0.53
54:DX:62:LYS:CD	54:DX:68:ARG:HD2	2.39	0.53
56:DZ:8:TYR:O	56:DZ:37:VAL:HB	2.09	0.53
1:AA:1023:G:H2'	1:AA:1024:G:H5'	1.91	0.53
1:AA:1108:G:OP2	3:AC:174:PRO:HA	2.09	0.53
1:AA:1207:G:O2'	1:AA:1208:C:H5'	2.09	0.53
1:AA:1253:G:H2'	1:AA:1254:C:H6	1.74	0.53
1:AA:1298:C:C6	7:AG:114:ARG:NH1	2.76	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:340:U:H2'	1:AA:341:C:C6	2.43	0.53
1:AA:702:A:H3'	1:AA:703:G:C5'	2.38	0.53
1:AA:745:C:H2'	1:AA:746:A:C8	2.44	0.53
1:AA:979:C:OP1	1:AA:1223:C:N4	2.42	0.53
1:AA:985:C:H2'	1:AA:986:A:H8	1.73	0.53
2:AB:80:ILE:HD13	2:AB:212:GLN:HA	1.91	0.53
3:AC:57:ILE:HG23	3:AC:64:VAL:CG1	2.39	0.53
3:AC:84:ILE:O	3:AC:88:ARG:HG3	2.09	0.53
5:AE:145:LYS:HG2	5:AE:149:GLU:OE1	2.09	0.53
5:AE:78:HIS:HD2	8:AH:104:ARG:HE	1.56	0.53
11:AK:122:LYS:O	11:AK:126:ARG:CB	2.57	0.53
14:AN:34:TYR:C	14:AN:36:PHE:N	2.62	0.53
17:AQ:71:PHE:N	17:AQ:71:PHE:HD2	2.07	0.53
17:AQ:71:PHE:N	17:AQ:71:PHE:CD2	2.76	0.53
17:AQ:95:TYR:C	17:AQ:97:SER:N	2.61	0.53
18:AR:85:LEU:HG	18:AR:86:VAL:N	2.17	0.53
20:AT:56:MET:O	20:AT:60:GLU:HB2	2.07	0.53
25:AY:84:ARG:HE	25:AY:92:PRO:CD	2.19	0.53
27:B1:58:ILE:CD1	27:B1:59:THR:N	2.64	0.53
27:B1:86:SER:N	27:B1:87:PRO:CD	2.72	0.53
29:B3:21:ALA:O	29:B3:24:LYS:N	2.38	0.53
33:B7:30:VAL:CG1	33:B7:33:ARG:HH22	2.22	0.53
35:BA:1407:C:O2	35:BA:1407:C:H2'	2.08	0.53
35:BA:237:C:H2'	35:BA:238:C:C6	2.42	0.53
35:BA:2380:C:H2'	35:BA:2381:C:H6	1.73	0.53
35:BA:251:A:H5''	46:BP:51:PHE:HZ	1.66	0.53
35:BA:2574:G:H2'	35:BA:2575:C:C6	2.44	0.53
35:BA:2658:C:H41	35:BA:2664:G:N2	2.06	0.53
35:BA:272(B):G:H2'	35:BA:272(C):G:H8	1.73	0.53
35:BA:2790:A:H2'	35:BA:2790:A:N3	2.23	0.53
35:BA:558:G:OP1	44:BN:111:PRO:HD2	2.09	0.53
35:BA:845:G:H8	35:BA:845:G:OP2	1.92	0.53
36:BB:28:C:N4	36:BB:56:G:H1	2.07	0.53
40:BF:45:ARG:HG3	40:BF:46:ARG:H	1.74	0.53
41:BG:55:LYS:C	41:BG:57:ALA:N	2.62	0.53
41:BG:67:LYS:O	41:BG:92:VAL:HG23	2.09	0.53
43:BI:114:LEU:O	43:BI:115:ALA:HB3	2.09	0.53
43:BI:89:TYR:O	43:BI:121:LYS:HE2	2.08	0.53
43:BI:123:LEU:HG	43:BI:142:VAL:HB	1.89	0.53
43:BI:5:LEU:O	43:BI:6:LEU:HG	2.08	0.53
44:BN:3:THR:C	44:BN:5:VAL:H	2.12	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:87:ILE:HD12	45:BO:92:GLU:N	2.24	0.53
46:BP:96:THR:HB	46:BP:97:PRO:CD	2.38	0.53
47:BQ:134:ARG:CG	47:BQ:135:ASP:H	2.13	0.53
47:BQ:28:ALA:HB1	47:BQ:29:PHE:CD1	2.44	0.53
48:BR:35:THR:HG23	48:BR:112:ALA:O	2.09	0.53
49:BS:92:TYR:HB3	49:BS:97:ARG:HH11	1.74	0.53
50:BT:81:PRO:O	50:BT:82:LEU:HD12	2.09	0.53
52:BV:19:LYS:C	52:BV:20:LEU:HD12	2.29	0.53
52:BV:89:GLN:HA	52:BV:89:GLN:NE2	2.24	0.53
53:BW:18:ARG:HG2	53:BW:18:ARG:HH11	1.74	0.53
35:BA:1312:U:OP2	54:BX:62:LYS:HE2	2.08	0.53
1:CA:1155:G:C2'	1:CA:1156:G:H5'	2.39	0.53
1:CA:1238:A:H62	1:CA:1301:U:H3	1.56	0.53
2:CB:144:ARG:HG3	2:CB:145:LEU:N	2.24	0.53
2:CB:74:LYS:CG	2:CB:77:ALA:HB3	2.39	0.53
3:CC:84:ILE:O	3:CC:88:ARG:HG3	2.09	0.53
4:CD:196:LEU:HD12	4:CD:196:LEU:H	1.74	0.53
4:CD:59:ARG:HH11	4:CD:59:ARG:HG2	1.73	0.53
5:CE:80:ILE:CG1	5:CE:91:LEU:HB2	2.38	0.53
8:CH:27:PRO:HG3	8:CH:58:TYR:CE2	2.44	0.53
8:CH:7:ALA:O	8:CH:8:ASP:C	2.47	0.53
14:CN:42:ILE:O	14:CN:43:CYS:C	2.47	0.53
14:CN:44:LEU:O	14:CN:44:LEU:HD12	2.09	0.53
16:CP:48:TRP:HE3	16:CP:49:LEU:N	2.06	0.53
16:CP:49:LEU:HD22	16:CP:73:LEU:HD13	1.90	0.53
19:CS:16:LEU:HB3	19:CS:20:LEU:HD11	1.91	0.53
19:CS:18:LYS:HA	19:CS:21:GLU:HG2	1.90	0.53
32:D6:10:LEU:H	32:D6:10:LEU:HD22	1.73	0.53
35:DA:1014:U:H2'	35:DA:1015:G:O4'	2.08	0.53
35:DA:1142(A):A:C5	35:DA:1144:G:C5	2.96	0.53
35:DA:1238:G:H2'	35:DA:1239:G:H8	1.74	0.53
35:DA:1543:C:H2'	35:DA:1543:C:O2	2.08	0.53
35:DA:1754:C:H2'	35:DA:1755:A:O4'	2.09	0.53
35:DA:2119:A:H3'	35:DA:2120:G:H5''	1.91	0.53
35:DA:2351:G:HO2'	35:DA:2352:A:H8	1.57	0.53
35:DA:2843:G:H1	35:DA:2874:C:N4	2.06	0.53
35:DA:543:C:H6	35:DA:547:A:C8	2.26	0.53
35:DA:684:G:H22	35:DA:787:U:H2'	1.72	0.53
38:DD:125:ILE:CD1	38:DD:125:ILE:N	2.72	0.53
38:DD:182:LEU:O	38:DD:271:ILE:HD12	2.08	0.53
44:DN:76:SER:O	44:DN:77:GLY:C	2.47	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:97:ARG:HG3	45:DO:97:ARG:HH11	1.74	0.53
35:DA:2415:G:C3'	46:DP:66:GLY:HA3	2.37	0.53
49:DS:11:LYS:HD3	49:DS:11:LYS:N	2.24	0.53
49:DS:26:LEU:HD23	49:DS:28:VAL:HG22	1.91	0.53
50:DT:53:ARG:CG	50:DT:53:ARG:NH1	2.72	0.53
51:DU:68:ALA:O	51:DU:71:GLN:N	2.41	0.53
51:DU:90:VAL:CG1	52:DV:39:LEU:HB3	2.35	0.53
54:DX:12:VAL:HG11	54:DX:27:THR:CG2	2.37	0.53
54:DX:39:ILE:C	54:DX:39:ILE:HD12	2.29	0.53
54:DX:83:VAL:O	54:DX:84:ALA:CB	2.56	0.53
1:AA:102:G:H2'	1:AA:103:C:C6	2.44	0.53
1:AA:1067:A:O5'	1:AA:1067:A:H8	1.91	0.53
1:AA:1155:G:O2'	1:AA:1156:G:H5'	2.09	0.53
1:AA:1217:C:H2'	1:AA:1218:C:C6	2.44	0.53
1:AA:1256:A:N6	1:AA:1278:U:H1'	2.24	0.53
1:AA:544:G:H2'	1:AA:545:C:H6	1.74	0.53
1:AA:678:U:H2'	1:AA:679:C:H6	1.66	0.53
3:AC:5:ILE:HD13	3:AC:5:ILE:O	2.08	0.53
6:AF:19:LEU:HD23	6:AF:19:LEU:C	2.28	0.53
7:AG:47:CYS:HB3	7:AG:58:PRO:HG3	1.90	0.53
8:AH:29:SER:O	8:AH:32:LYS:HB2	2.08	0.53
1:AA:599:C:O3'	8:AH:96:GLY:HA2	2.09	0.53
12:AL:89:ARG:NH1	12:AL:90:VAL:N	2.57	0.53
15:AO:17:ARG:NH1	15:AO:17:ARG:HG3	2.23	0.53
16:AP:48:TRP:CE3	16:AP:49:LEU:HB3	2.44	0.53
27:B1:66:HIS:C	27:B1:68:PRO:HD2	2.29	0.53
27:B1:64:ALA:C	27:B1:67:ILE:HG13	2.27	0.53
32:B6:36:LEU:HD13	32:B6:50:ARG:HH12	1.71	0.53
34:B8:30:ARG:HE	46:BP:62:LEU:CB	2.22	0.53
35:BA:1215:G:H2'	35:BA:1216:G:C8	2.44	0.53
35:BA:1430:C:H2'	35:BA:1431:U:C6	2.44	0.53
35:BA:1485:G:H2'	35:BA:1486:A:H8	1.73	0.53
35:BA:1630:G:H2'	35:BA:1631:C:H6	1.74	0.53
35:BA:1856:G:H2'	35:BA:1857:G:H5'	1.91	0.53
35:BA:2639:A:C3'	35:BA:2640:G:C5'	2.86	0.53
35:BA:2768:C:O2'	35:BA:2769:C:H5'	2.09	0.53
35:BA:35:G:O2'	35:BA:36:G:H5'	2.08	0.53
35:BA:807:U:O2'	35:BA:808:G:H5'	2.09	0.53
36:BB:56:G:H4'	36:BB:57:A:C8	2.43	0.53
38:BD:133:LEU:O	38:BD:134:ARG:C	2.47	0.53
38:BD:193:VAL:O	38:BD:193:VAL:HG13	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:24:THR:OG1	39:BE:188:VAL:HG11	2.08	0.53
40:BF:68:LYS:O	40:BF:69:HIS:HB2	2.09	0.53
41:BG:94:LEU:HD13	41:BG:102:PHE:CD1	2.44	0.53
43:BI:10:GLU:CD	43:BI:11:ASN:N	2.63	0.53
35:BA:252:G:OP2	46:BP:50:ARG:NH1	2.42	0.53
49:BS:83:LYS:O	49:BS:85:VAL:HG22	2.09	0.53
49:BS:92:TYR:O	49:BS:93:LYS:HB2	2.07	0.53
50:BT:28:VAL:O	50:BT:29:ARG:HD3	2.09	0.53
50:BT:58:ASN:ND2	50:BT:58:ASN:N	2.56	0.53
51:BU:65:ILE:HD12	51:BU:65:ILE:N	2.24	0.53
51:BU:95:LEU:HA	51:BU:97:ASP:OD1	2.09	0.53
54:BX:57:LEU:HD12	54:BX:76:ARG:CD	2.39	0.53
56:BZ:103:ARG:CB	56:BZ:138:GLU:HA	2.32	0.53
56:BZ:149:SER:HB2	56:BZ:172:ALA:O	2.09	0.53
47:BQ:141:GLN:HG2	56:BZ:72:ARG:HE	1.73	0.53
1:CA:231:G:O2'	1:CA:232:G:H5'	2.09	0.53
1:CA:340:U:H2'	1:CA:341:C:C6	2.41	0.53
1:CA:940:C:H2'	1:CA:941:G:H8	1.74	0.53
1:CA:1096:C:H5''	2:CB:137:ARG:NH2	2.24	0.53
2:CB:89:GLY:O	2:CB:154:LEU:HD13	2.08	0.53
2:CB:193:ASP:OD2	2:CB:196:LEU:HD21	2.08	0.53
3:CC:70:VAL:O	3:CC:105:GLU:HA	2.09	0.53
5:CE:57:LYS:HE2	5:CE:61:TYR:CE2	2.42	0.53
9:CI:28:VAL:HG13	9:CI:63:ILE:C	2.29	0.53
12:CL:89:ARG:C	12:CL:89:ARG:HD3	2.29	0.53
13:CM:69:GLU:CB	13:CM:72:ALA:HB3	2.38	0.53
14:CN:34:TYR:C	14:CN:36:PHE:N	2.62	0.53
25:CY:173:ASP:O	25:CY:175:LEU:N	2.42	0.53
28:D2:15:LYS:O	28:D2:19:VAL:HG23	2.09	0.53
29:D3:40:THR:HG23	29:D3:43:ILE:H	1.72	0.53
29:D3:7:LYS:C	29:D3:54:VAL:HG13	2.28	0.53
32:D6:47:THR:HG22	32:D6:48:VAL:N	2.24	0.53
35:DA:1024:G:C8	35:DA:1025:G:H2'	2.43	0.53
35:DA:443:A:H1'	35:DA:1201:C:O4'	2.08	0.53
35:DA:1518:U:H2'	35:DA:1519:G:C8	2.43	0.53
35:DA:1665:A:H1'	45:DO:1:MET:HE3	1.91	0.53
35:DA:210:C:O2'	35:DA:211:A:H5'	2.09	0.53
35:DA:2260:C:H2'	35:DA:2261:C:H6	1.74	0.53
35:DA:2380:C:H2'	35:DA:2381:C:H6	1.74	0.53
35:DA:2528:U:H2'	35:DA:2530:A:O5'	2.09	0.53
35:DA:300:A:H5''	55:DY:97:ARG:HH12	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:524:U:H4'	35:DA:555:U:H4'	1.91	0.53
35:DA:71:A:H5'	35:DA:71:A:H8	1.73	0.53
39:DE:55:ASN:CG	39:DE:75:VAL:HG13	2.29	0.53
39:DE:77:ILE:HG23	39:DE:78:LEU:H	1.73	0.53
40:DF:9:ILE:O	40:DF:128:ALA:HB2	2.09	0.53
41:DG:115:ARG:HH12	41:DG:136:ARG:HD2	1.74	0.53
44:DN:74:ARG:NH2	44:DN:101:HIS:O	2.39	0.53
44:DN:42:TRP:N	51:DU:64:ARG:HH12	2.06	0.53
35:DA:1952:A:C5	45:DO:22:ILE:HD12	2.44	0.53
46:DP:122:PRO:HA	46:DP:141:ALA:O	2.08	0.53
46:DP:74:GLU:OE2	46:DP:74:GLU:HA	2.08	0.53
46:DP:96:THR:O	46:DP:99:LEU:HB3	2.09	0.53
50:DT:63:VAL:O	50:DT:73:GLU:HA	2.09	0.53
51:DU:24:TYR:HB2	51:DU:29:SER:CB	2.39	0.53
52:DV:19:LYS:C	52:DV:20:LEU:HD12	2.29	0.53
52:DV:32:THR:CG2	52:DV:33:VAL:H	2.18	0.53
52:DV:36:PRO:HG2	52:DV:60:GLU:OE1	2.09	0.53
52:DV:37:VAL:HG11	52:DV:53:GLU:OE2	2.09	0.53
56:DZ:48:PHE:HA	56:DZ:52:SER:H	1.73	0.53
56:DZ:55:HIS:CE1	56:DZ:135:GLU:HA	2.44	0.53
1:AA:1444:C:H2'	1:AA:1445:C:C6	2.43	0.53
1:AA:164:U:O2'	1:AA:165:C:H5'	2.08	0.53
1:AA:788:U:O2'	1:AA:789:U:H5'	2.08	0.53
2:AB:72:GLY:HA3	2:AB:165:VAL:CG2	2.38	0.53
6:AF:60:PHE:O	6:AF:61:LEU:HD12	2.09	0.53
7:AG:47:CYS:O	7:AG:58:PRO:HG3	2.09	0.53
9:AI:105:ASP:C	9:AI:107:ARG:H	2.11	0.53
11:AK:43:SER:CA	11:AK:47:VAL:HG21	2.37	0.53
13:AM:2:ALA:O	13:AM:9:ILE:HG13	2.09	0.53
19:AS:9:VAL:HG12	19:AS:9:VAL:O	2.09	0.53
20:AT:63:ILE:HD12	20:AT:81:LYS:CG	2.39	0.53
27:B1:20:ARG:CZ	27:B1:41:ARG:HE	2.20	0.53
27:B1:54:ALA:HB2	27:B1:57:GLU:OE1	2.09	0.53
31:B5:2:ALA:HB2	35:BA:2014:A:HO2'	1.72	0.53
31:B5:7:PRO:HA	35:BA:2615:U:C2	2.44	0.53
35:BA:1297:C:H2'	35:BA:1298:C:H6	1.74	0.53
35:BA:1348:G:C3'	35:BA:1349:A:H5''	2.39	0.53
35:BA:1487:G:N2	35:BA:1488:G:H1'	2.24	0.53
35:BA:1721:G:H2'	35:BA:1722:A:H2'	1.90	0.53
35:BA:1835:G:C5'	35:BA:1836:C:OP2	2.56	0.53
35:BA:2248:C:H2'	35:BA:2249:U:H5'	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2314:C:O2'	35:BA:2315:G:H5'	2.09	0.53
35:BA:235:U:H2'	35:BA:236:C:H6	1.73	0.53
35:BA:2383:G:O2'	35:BA:2384:G:H5'	2.09	0.53
35:BA:2685:G:C2	35:BA:2686:G:N7	2.76	0.53
35:BA:2821:A:OP2	48:BR:2:ARG:NH2	2.41	0.53
35:BA:631:A:H4'	46:BP:65:ARG:HA	1.91	0.53
35:BA:858:U:O2	35:BA:2268:A:H2'	2.09	0.53
37:BC:203:GLY:O	37:BC:204:ALA:HB2	2.09	0.53
37:BC:49:ILE:C	37:BC:51:PRO:HD3	2.29	0.53
39:BE:171:GLU:O	39:BE:173:VAL:HG23	2.08	0.53
39:BE:49:LEU:HD23	39:BE:81:ILE:HG12	1.91	0.53
41:BG:6:ALA:O	41:BG:10:LYS:N	2.31	0.53
42:BH:164:TYR:O	42:BH:165:ALA:HB2	2.08	0.53
43:BI:17:GLN:HG2	43:BI:18:VAL:N	2.21	0.53
46:BP:65:ARG:HH11	46:BP:65:ARG:HG3	1.74	0.53
47:BQ:26:TYR:O	47:BQ:27:VAL:C	2.48	0.53
49:BS:42:ASP:O	49:BS:44:LYS:N	2.42	0.53
49:BS:53:SER:OG	49:BS:54:LEU:N	2.42	0.53
50:BT:28:VAL:CG2	50:BT:46:GLU:HA	2.39	0.53
50:BT:80:SER:O	50:BT:82:LEU:N	2.41	0.53
51:BU:3:ARG:NH1	51:BU:3:ARG:CG	2.71	0.53
52:BV:36:PRO:HG2	52:BV:60:GLU:OE1	2.07	0.53
52:BV:5:VAL:HG22	52:BV:6:LYS:N	2.24	0.53
53:BW:14:PRO:HG3	53:BW:101:SER:OG	2.09	0.53
35:BA:1614:A:N1	53:BW:87:PRO:HB3	2.23	0.53
54:BX:52:VAL:O	54:BX:53:LYS:HB3	2.09	0.53
55:BY:29:GLU:N	55:BY:29:GLU:OE1	2.42	0.53
56:BZ:150:LEU:CD2	56:BZ:150:LEU:H	2.11	0.53
56:BZ:166:SER:HB2	56:BZ:168:GLU:H	1.74	0.53
1:CA:1090:U:O2'	1:CA:1091:U:H5'	2.09	0.53
1:CA:1116:C:H3'	1:CA:1117:G:H5''	1.91	0.53
1:CA:176:C:H2'	1:CA:177:C:C6	2.44	0.53
1:CA:223:U:H2'	1:CA:224:C:H6	1.74	0.53
1:CA:527:G:O2'	1:CA:528:C:H5'	2.09	0.53
1:CA:76:C:H42	1:CA:93:G:H1	1.57	0.53
1:CA:788:U:C5	1:CA:789:U:H5	2.26	0.53
1:CA:792:A:H4'	1:CA:793:U:O5'	2.08	0.53
1:CA:969:A:H2'	1:CA:970:C:O4'	2.09	0.53
5:CE:145:LYS:HG2	5:CE:149:GLU:OE1	2.08	0.53
7:CG:50:ILE:HG13	7:CG:58:PRO:HB3	1.91	0.53
11:CK:33:THR:C	11:CK:40:ILE:HG12	2.30	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:119:LYS:C	12:CL:121:GLY:H	2.11	0.53
13:CM:22:ILE:HG21	13:CM:25:ILE:HD12	1.90	0.53
14:CN:39:LEU:HD11	14:CN:47:LEU:HD12	1.91	0.53
15:CO:6:GLU:CD	15:CO:6:GLU:H	2.10	0.53
23:CW:39:A:H2'	23:CW:40:C:H5'	1.90	0.53
25:CY:164:ILE:HD12	25:CY:164:ILE:N	2.22	0.53
35:DA:1188:U:H2'	35:DA:1189:A:H5'	1.91	0.53
35:DA:1799:G:O3'	38:DD:264:LYS:NZ	2.42	0.53
35:DA:1960:A:O2'	35:DA:1961:C:H5'	2.09	0.53
35:DA:2114:A:N3	35:DA:2114:A:H2'	2.23	0.53
35:DA:2711:A:OP1	35:DA:2712(A):A:P	2.67	0.53
35:DA:443:A:P	35:DA:614(B):G:H22	2.32	0.53
36:DB:104:U:H2'	36:DB:105:A:H8	1.74	0.53
36:DB:78:A:O2'	36:DB:79:C:H5'	2.09	0.53
36:DB:95:C:O2'	36:DB:96:U:H5'	2.08	0.53
35:DA:1569:A:O2'	38:DD:38:LYS:HE2	2.09	0.53
39:DE:34:VAL:O	39:DE:35:GLN:HB2	2.09	0.53
40:DF:31:HIS:HB2	46:DP:13:ASN:OD1	2.09	0.53
42:DH:89:ILE:CG1	42:DH:90:LYS:H	2.22	0.53
45:DO:24:VAL:HG21	45:DO:32:TYR:O	2.09	0.53
46:DP:124:LYS:HA	46:DP:143:GLY:HA3	1.91	0.53
48:DR:38:VAL:HB	48:DR:39:PRO:CD	2.29	0.53
48:DR:44:LEU:HD13	48:DR:48:VAL:HG23	1.91	0.53
48:DR:73:VAL:HG23	48:DR:74:LYS:HD2	1.90	0.53
53:DW:40:ASN:O	53:DW:41:LYS:HG2	2.09	0.53
53:DW:51:LEU:O	53:DW:51:LEU:HD22	2.09	0.53
55:DY:37:VAL:CG2	55:DY:67:LEU:HG	2.39	0.53
56:DZ:26:GLY:O	56:DZ:37:VAL:N	2.42	0.53
1:AA:1053:G:C3'	1:AA:1054:C:H5'	2.38	0.52
1:AA:1404:C:H6	1:AA:1404:C:O5'	1.92	0.52
1:AA:358:U:O2'	1:AA:359:U:H5'	2.09	0.52
5:AE:57:LYS:HE2	5:AE:61:TYR:CE2	2.44	0.52
6:AF:75:LEU:HD22	6:AF:79:LEU:HD11	1.91	0.52
8:AH:6:ILE:CG2	8:AH:10:LEU:HD11	2.38	0.52
8:AH:124:ALA:O	8:AH:128:GLY:N	2.40	0.52
12:AL:8:ASN:O	12:AL:11:VAL:HB	2.09	0.52
13:AM:82:MET:HB3	13:AM:93:ARG:NH1	2.24	0.52
15:AO:67:LEU:HD22	15:AO:78:TYR:CE1	2.44	0.52
16:AP:1:MET:SD	16:AP:1:MET:N	2.74	0.52
17:AQ:86:GLU:O	17:AQ:88:TYR:N	2.41	0.52
19:AS:16:LEU:HB3	19:AS:20:LEU:HD11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:91:ARG:HD3	19:AS:81:ARG:NH2	2.24	0.52
22:AV:40:C:H2'	22:AV:41:C:H6	1.73	0.52
25:AY:37:LEU:HD12	25:AY:38:LEU:HG	1.91	0.52
27:B1:20:ARG:HH12	27:B1:41:ARG:NE	2.06	0.52
32:B6:9:LEU:HD23	32:B6:10:LEU:O	2.09	0.52
35:BA:1038:C:H42	35:BA:1117:G:H1	1.57	0.52
35:BA:1175:U:H4'	35:BA:1176:G:C3'	2.38	0.52
35:BA:818:G:N2	35:BA:1190:G:C6	2.77	0.52
35:BA:188:G:H1'	35:BA:1365:A:N1	2.23	0.52
35:BA:2881:C:C2	35:BA:2882:A:C8	2.97	0.52
35:BA:66:C:H2'	35:BA:67:U:C5'	2.39	0.52
37:BC:44:HIS:HD2	37:BC:175:VAL:N	2.06	0.52
37:BC:99:ILE:HG23	37:BC:103:ILE:CB	2.38	0.52
38:BD:132:PRO:HD3	38:BD:190:TYR:CZ	2.44	0.52
38:BD:83:GLU:HB2	38:BD:92:ILE:HD11	1.91	0.52
40:BF:34:TRP:O	40:BF:37:VAL:N	2.41	0.52
41:BG:107:LEU:HD23	41:BG:111:LEU:HD12	1.90	0.52
42:BH:128:PRO:HG2	42:BH:129:THR:CG2	2.35	0.52
43:BI:29:TYR:HD1	43:BI:33:ARG:HE	1.57	0.52
45:BO:34:THR:O	45:BO:37:ASP:OD2	2.27	0.52
45:BO:63:VAL:HG22	45:BO:84:ALA:CA	2.25	0.52
46:BP:70:GLN:HG3	46:BP:71:VAL:N	2.21	0.52
46:BP:97:PRO:C	46:BP:99:LEU:H	2.11	0.52
50:BT:22:PHE:O	50:BT:23:ARG:HB3	2.09	0.52
50:BT:53:ARG:HH11	50:BT:53:ARG:CG	2.21	0.52
51:BU:3:ARG:HG2	51:BU:3:ARG:NH1	2.19	0.52
51:BU:78:THR:O	51:BU:80:ILE:N	2.42	0.52
52:BV:19:LYS:HG3	52:BV:20:LEU:H	1.74	0.52
52:BV:96:ILE:HG23	52:BV:97:LYS:H	1.73	0.52
53:BW:66:GLU:HG2	53:BW:66:GLU:O	2.08	0.52
55:BY:47:LYS:HG3	55:BY:60:PHE:CE2	2.44	0.52
56:BZ:80:ARG:O	56:BZ:82:ARG:N	2.42	0.52
1:CA:1053:G:C6	1:CA:1199:U:H2'	2.42	0.52
1:CA:1496:C:H2'	1:CA:1497:G:C8	2.43	0.52
2:CB:158:LEU:H	2:CB:158:LEU:HD12	1.72	0.52
3:CC:57:ILE:HG23	3:CC:64:VAL:CG1	2.39	0.52
3:CC:73:PRO:O	3:CC:75:VAL:N	2.42	0.52
11:CK:125:PHE:CD1	11:CK:125:PHE:N	2.77	0.52
11:CK:20:TYR:C	11:CK:21:ILE:HD12	2.29	0.52
25:CY:114:LEU:O	25:CY:117:ALA:HB3	2.09	0.52
27:D1:9:GLY:N	27:D1:48:LYS:HZ3	1.97	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:D1:76:ARG:C	27:D1:78:LYS:HZ2	2.13	0.52
34:D8:29:LYS:O	34:D8:30:ARG:C	2.47	0.52
34:D8:30:ARG:HE	46:DP:62:LEU:CB	2.21	0.52
35:DA:818:G:OP2	35:DA:1187:G:O6	2.27	0.52
35:DA:1666:G:O3'	45:DO:6:THR:HG23	2.09	0.52
35:DA:154(A):C:C5	35:DA:171:G:N1	2.75	0.52
35:DA:1744:C:C2'	35:DA:1745:C:H5'	2.38	0.52
35:DA:1824:G:H2'	35:DA:1825:A:H8	1.73	0.52
35:DA:1836:C:O2'	35:DA:1837:C:H5'	2.09	0.52
35:DA:224:G:O2'	35:DA:225:A:H5'	2.09	0.52
35:DA:247:G:H4'	35:DA:386:G:C5	2.45	0.52
35:DA:26:G:H1'	35:DA:515:A:N6	2.15	0.52
35:DA:27:G:H1'	35:DA:513:A:H62	1.75	0.52
35:DA:463:G:N1	35:DA:467:G:C6	2.77	0.52
35:DA:516:C:O2'	35:DA:517:C:H5'	2.09	0.52
35:DA:587:C:C6	35:DA:671:C:H1'	2.44	0.52
35:DA:668:G:C2	35:DA:670:A:C6	2.97	0.52
36:DB:55:U:H4'	41:DG:27:ASN:HD21	1.72	0.52
40:DF:68:LYS:O	40:DF:69:HIS:CB	2.57	0.52
40:DF:9:ILE:HG12	40:DF:14:PRO:C	2.29	0.52
41:DG:125:PHE:CE2	41:DG:173:LEU:HD12	2.43	0.52
43:DI:130:TYR:CB	43:DI:136:VAL:HG13	2.38	0.52
43:DI:83:ALA:HB2	43:DI:88:ILE:HG23	1.91	0.52
45:DO:13:ASN:HD22	45:DO:97:ARG:HB2	1.74	0.52
46:DP:71:VAL:HG22	46:DP:72:PRO:N	2.24	0.52
35:DA:2009:G:H1'	48:DR:107:ASP:C	2.29	0.52
49:DS:65:VAL:HG12	49:DS:69:VAL:HB	1.91	0.52
50:DT:36:GLU:HB3	50:DT:38:ASN:OD1	2.09	0.52
53:DW:29:LEU:HD23	53:DW:29:LEU:C	2.30	0.52
53:DW:73:ALA:HB3	53:DW:106:ILE:CD1	2.26	0.52
55:DY:87:LYS:C	55:DY:88:LYS:HD2	2.29	0.52
56:DZ:144:LEU:HD21	56:DZ:150:LEU:CD1	2.38	0.52
1:AA:1366:C:H2'	1:AA:1367:C:H6	1.70	0.52
1:AA:139:G:O2'	1:AA:140:A:H5'	2.09	0.52
1:AA:1418:A:C2	1:AA:1483:A:C2	2.97	0.52
1:AA:184:G:O2'	1:AA:185:A:H5'	2.09	0.52
1:AA:644:G:O2'	1:AA:645:C:H5'	2.09	0.52
2:AB:115:LEU:HD21	2:AB:153:ARG:HE	1.74	0.52
2:AB:36:ARG:N	2:AB:41:ILE:HD13	2.23	0.52
4:AD:65:ARG:HH11	4:AD:72:GLU:CA	2.22	0.52
5:AE:15:ARG:HG2	5:AE:26:PHE:HD2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:105:VAL:O	7:AG:108:ALA:HB3	2.10	0.52
9:AI:10:ARG:O	9:AI:11:LYS:HB3	2.09	0.52
9:AI:7:THR:N	9:AI:83:ARG:HD2	2.23	0.52
10:AJ:62:HIS:H	10:AJ:62:HIS:HD2	1.54	0.52
11:AK:21:ILE:CG1	11:AK:84:VAL:HG12	2.39	0.52
1:AA:624:C:H4'	16:AP:10:GLY:HA2	1.91	0.52
16:AP:1:MET:SD	16:AP:3:LYS:HE3	2.49	0.52
16:AP:72:ARG:C	16:AP:74:LEU:N	2.61	0.52
20:AT:48:LYS:HB2	20:AT:52:ALA:HB2	1.91	0.52
26:B0:27:GLU:N	26:B0:69:PHE:CE1	2.74	0.52
28:B2:51:ARG:HD3	28:B2:51:ARG:O	2.09	0.52
29:B3:7:LYS:C	29:B3:54:VAL:HG13	2.29	0.52
35:BA:109:G:H2'	35:BA:110:G:H8	1.75	0.52
35:BA:1297:C:H2'	35:BA:1298:C:C6	2.43	0.52
35:BA:1434:A:H2'	35:BA:1435:G:H8	1.74	0.52
35:BA:1518:U:H2'	35:BA:1519:G:C8	2.44	0.52
35:BA:2865:U:H3'	35:BA:2866:U:O2	2.09	0.52
35:BA:32:C:H42	35:BA:473:G:H1	1.57	0.52
35:BA:793:A:OP2	35:BA:2071:A:O2'	2.25	0.52
36:BB:87:G:H2'	36:BB:88:C:H5''	1.90	0.52
38:BD:130:ALA:HB2	38:BD:192:THR:HB	1.91	0.52
35:BA:1827:C:OP2	38:BD:222:ARG:NH1	2.42	0.52
39:BE:55:ASN:CG	39:BE:75:VAL:HG13	2.29	0.52
44:BN:56:ASN:HA	44:BN:125:GLY:N	2.23	0.52
46:BP:70:GLN:HG3	46:BP:71:VAL:HG12	1.91	0.52
47:BQ:132:VAL:HG12	47:BQ:133:ARG:N	2.25	0.52
49:BS:83:LYS:O	49:BS:85:VAL:HG13	2.09	0.52
50:BT:25:GLY:HA2	50:BT:92:GLY:HA3	1.91	0.52
44:BN:40:PRO:CB	51:BU:64:ARG:HH22	2.22	0.52
53:BW:74:ALA:O	53:BW:75:TYR:HB3	2.09	0.52
54:BX:82:GLN:OE1	54:BX:83:VAL:N	2.38	0.52
47:BQ:141:GLN:HG2	56:BZ:72:ARG:NE	2.24	0.52
56:BZ:99:TYR:O	56:BZ:100:VAL:HB	2.08	0.52
1:CA:1051:C:O2'	1:CA:1052:U:H5'	2.09	0.52
1:CA:1096:C:O2'	1:CA:1097:C:H5'	2.09	0.52
1:CA:1234:C:H4'	1:CA:1364:U:H1'	1.91	0.52
1:CA:1287:A:C6	1:CA:1288:A:N6	2.76	0.52
1:CA:1442(A):G:N2	50:DT:119:LYS:HB2	2.22	0.52
1:CA:745:C:H2'	1:CA:746:A:C8	2.44	0.52
1:CA:833:U:H2'	1:CA:834:C:C6	2.45	0.52
1:CA:967:C:H2'	1:CA:968:A:C8	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:76:VAL:HG23	3:CC:77:ILE:HG13	1.90	0.52
4:CD:96:LEU:O	4:CD:98:GLU:N	2.43	0.52
10:CJ:85:LEU:O	10:CJ:87:THR:N	2.43	0.52
10:CJ:92:THR:HG23	10:CJ:93:GLY:N	2.24	0.52
11:CK:87:THR:HA	11:CK:91:ARG:CG	2.39	0.52
12:CL:60:LEU:HD22	12:CL:60:LEU:N	2.24	0.52
16:CP:19:ILE:HB	16:CP:37:GLY:CA	2.39	0.52
17:CQ:10:VAL:HG21	17:CQ:55:ASP:HB2	1.90	0.52
19:CS:27:GLU:O	19:CS:28:LYS:HD2	2.07	0.52
19:CS:53:ASN:N	19:CS:53:ASN:HD22	2.07	0.52
19:CS:62:ILE:HD12	19:CS:63:THR:N	2.24	0.52
20:CT:48:LYS:HB2	20:CT:52:ALA:HB2	1.91	0.52
25:CY:28:LEU:O	25:CY:37:LEU:HD13	2.09	0.52
25:CY:78:ALA:HA	25:CY:81:LYS:HB2	1.91	0.52
32:D6:20:ASN:HD22	32:D6:21:TYR:H	1.53	0.52
35:DA:1197:G:H2'	35:DA:1198:U:C6	2.44	0.52
35:DA:1210:A:O2'	35:DA:1211:U:OP2	2.26	0.52
35:DA:1341:U:OP1	35:DA:1397:U:N3	2.41	0.52
35:DA:1434:A:H2'	35:DA:1435:G:H8	1.68	0.52
35:DA:1419:A:H62	35:DA:1578:U:H3	1.58	0.52
35:DA:2290:G:H2'	35:DA:2291:U:C6	2.44	0.52
35:DA:262:A:H2'	35:DA:263:C:O4'	2.08	0.52
35:DA:535:C:C2'	35:DA:536:A:H5'	2.39	0.52
35:DA:680:G:H2'	35:DA:681:G:C8	2.44	0.52
35:DA:874:G:H2'	35:DA:875:G:C8	2.44	0.52
35:DA:910:A:H2'	35:DA:911:A:C8	2.44	0.52
37:DC:73:ARG:HA	37:DC:92:ASP:OD1	2.09	0.52
38:DD:9:TYR:C	38:DD:10:THR:HG22	2.29	0.52
40:DF:114:VAL:HG21	40:DF:202:PHE:CE2	2.44	0.52
41:DG:171:ALA:O	41:DG:174:GLU:N	2.42	0.52
41:DG:44:GLY:C	41:DG:46:ALA:N	2.63	0.52
42:DH:45:VAL:HG12	42:DH:45:VAL:O	2.08	0.52
43:DI:37:VAL:CG1	43:DI:38:LEU:N	2.72	0.52
45:DO:31:LYS:HD2	45:DO:32:TYR:CE1	2.44	0.52
45:DO:62:VAL:HG11	45:DO:65:THR:HG22	1.91	0.52
46:DP:13:ASN:HD22	46:DP:13:ASN:N	2.03	0.52
48:DR:87:TYR:O	48:DR:89:ASP:N	2.41	0.52
35:DA:993:G:OP1	51:DU:50:ARG:NH2	2.42	0.52
51:DU:61:TRP:O	51:DU:63:VAL:N	2.42	0.52
52:DV:5:VAL:HG22	52:DV:6:LYS:N	2.24	0.52
53:DW:48:ALA:O	53:DW:51:LEU:N	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:494:G:H21	53:DW:57:ASN:HD21	1.55	0.52
55:DY:18:GLY:C	55:DY:20:TYR:H	2.12	0.52
56:DZ:125:LEU:HD23	56:DZ:126:VAL:N	2.24	0.52
56:DZ:53:ILE:HG22	56:DZ:71:VAL:CG2	2.39	0.52
1:AA:1037:C:H2'	1:AA:1038:C:C6	2.44	0.52
1:AA:269:C:H2'	1:AA:270:A:H8	1.74	0.52
1:AA:364:A:H2'	1:AA:365:U:O2	2.09	0.52
1:AA:527:G:C2'	1:AA:528:C:H5'	2.39	0.52
1:AA:622:A:C8	1:AA:623:C:C5	2.97	0.52
12:AL:34:ARG:HB3	12:AL:61:THR:CG2	2.38	0.52
18:AR:59:SER:N	18:AR:62:GLU:OE1	2.42	0.52
19:AS:53:ASN:N	19:AS:53:ASN:ND2	2.55	0.52
21:AU:18:TYR:CD2	21:AU:24:ARG:HG3	2.44	0.52
27:B1:17:SER:O	27:B1:44:PRO:HD2	2.08	0.52
28:B2:47:ASN:HA	28:B2:50:ILE:O	2.09	0.52
31:B5:40:LYS:HZ2	31:B5:45:VAL:HA	1.72	0.52
34:B8:39:LYS:NZ	34:B8:43:GLN:HG3	2.24	0.52
34:B8:61:LEU:O	34:B8:64:TYR:HD1	1.93	0.52
35:BA:1006:C:N3	35:BA:1138:G:C2	2.77	0.52
35:BA:1197:G:H2'	35:BA:1198:U:C6	2.45	0.52
35:BA:149:A:C2	35:BA:150:C:C2	2.98	0.52
35:BA:1841:U:H2'	35:BA:1842:G:H8	1.72	0.52
35:BA:2171:A:HO2'	35:BA:2172:U:H6	1.55	0.52
35:BA:2262:U:H4'	35:BA:2328:A:H2	1.74	0.52
35:BA:2537:U:H2'	35:BA:2538:C:C6	2.43	0.52
35:BA:2809:A:C2	35:BA:2892:A:N3	2.77	0.52
35:BA:443:A:P	35:BA:614(B):G:H22	2.32	0.52
35:BA:589:C:O2'	35:BA:590:A:H5'	2.10	0.52
35:BA:705:A:O2'	35:BA:706:A:H5'	2.10	0.52
38:BD:201:HIS:C	38:BD:203:ASN:H	2.13	0.52
41:BG:116:ASP:CG	41:BG:117:PHE:H	2.12	0.52
41:BG:139:LEU:HD22	41:BG:146:TYR:HE1	1.74	0.52
36:BB:45:A:C8	41:BG:95:ARG:NE	2.78	0.52
42:BH:89:ILE:HD13	42:BH:89:ILE:H	1.75	0.52
43:BI:58:LEU:C	43:BI:58:LEU:HD23	2.30	0.52
46:BP:71:VAL:HG22	46:BP:72:PRO:N	2.24	0.52
35:BA:1652:A:H62	48:BR:11:ASN:ND2	2.07	0.52
50:BT:13:ARG:NH1	50:BT:15:VAL:HG12	2.24	0.52
51:BU:31:SER:O	51:BU:33:ARG:N	2.43	0.52
51:BU:74:LEU:CD1	51:BU:79:PHE:HB2	2.39	0.52
52:BV:37:VAL:HG11	52:BV:53:GLU:OE2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BV:93:GLU:HG2	52:BV:94:LEU:N	2.24	0.52
35:BA:494:G:H21	53:BW:57:ASN:HD21	1.55	0.52
54:BX:55:ASN:HB2	54:BX:77:LYS:CD	2.40	0.52
54:BX:73:ARG:HG3	54:BX:73:ARG:O	2.09	0.52
56:BZ:28:MET:HA	56:BZ:88:PHE:O	2.10	0.52
1:CA:1108:G:OP2	3:CC:174:PRO:HA	2.08	0.52
1:CA:129(A):G:H21	1:CA:189(F):U:H5''	1.74	0.52
1:CA:1320:C:H2'	1:CA:1321:C:O4'	2.09	0.52
1:CA:1367:C:OP1	9:CI:115:GLY:N	2.39	0.52
1:CA:237:C:O2'	1:CA:238:G:H5'	2.08	0.52
1:CA:429:U:C1'	1:CA:430:A:H5''	2.39	0.52
1:CA:475:G:H2'	1:CA:476:G:H8	1.74	0.52
2:CB:9:GLU:O	2:CB:13:ALA:HB2	2.10	0.52
2:CB:41:ILE:N	2:CB:41:ILE:HD12	2.24	0.52
4:CD:131:ARG:O	4:CD:132:ARG:C	2.46	0.52
4:CD:153:ARG:HB3	4:CD:153:ARG:HH11	1.73	0.52
4:CD:156:GLU:O	4:CD:159:ARG:HB2	2.10	0.52
5:CE:129:ILE:O	5:CE:130:ASN:C	2.48	0.52
7:CG:126:ASP:HB3	7:CG:132:GLY:CA	2.39	0.52
7:CG:49:ILE:HG22	7:CG:53:LYS:HG3	1.92	0.52
9:CI:11:LYS:C	9:CI:13:ALA:H	2.12	0.52
9:CI:65:VAL:CG2	9:CI:66:ARG:N	2.73	0.52
12:CL:24:VAL:O	12:CL:24:VAL:HG12	2.08	0.52
15:CO:9:GLN:HB3	15:CO:13:GLN:NE2	2.17	0.52
17:CQ:29:HIS:HA	17:CQ:36:ILE:HD11	1.90	0.52
25:CY:35:PRO:HA	25:CY:66:LEU:HD21	1.92	0.52
25:CY:73:GLN:O	25:CY:77:LYS:HE3	2.09	0.52
35:DA:1259:G:O2'	35:DA:1260:G:H5'	2.09	0.52
35:DA:1348:G:C3'	35:DA:1349:A:H5''	2.40	0.52
35:DA:1430:C:H2'	35:DA:1431:U:H6	1.72	0.52
35:DA:1469:A:O2'	35:DA:1470:G:H5'	2.08	0.52
35:DA:1860:G:H2'	35:DA:1861:G:H8	1.74	0.52
35:DA:189:G:C4	35:DA:205:G:N2	2.77	0.52
35:DA:2313:C:P	41:DG:71:THR:HG21	2.49	0.52
35:DA:271(H):G:HO2'	35:DA:271(I):G:H8	1.57	0.52
35:DA:2729:G:H2'	35:DA:2730:C:C6	2.44	0.52
35:DA:442:G:C4'	40:DF:46:ARG:HD3	2.39	0.52
35:DA:477:A:H2'	35:DA:478:A:C8	2.45	0.52
35:DA:579:G:H2'	35:DA:580:C:C6	2.45	0.52
35:DA:701:G:O2'	35:DA:702:G:H5'	2.09	0.52
35:DA:819:A:OP2	35:DA:1187:G:N2	2.31	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:927:G:O6	35:DA:928:G:C2	2.62	0.52
36:DB:114:C:O2'	49:DS:47:THR:HB	2.08	0.52
39:DE:50:GLY:CA	39:DE:74:PRO:HG3	2.40	0.52
40:DF:57:VAL:CG1	40:DF:58:ALA:N	2.72	0.52
30:D4:14:ILE:CB	41:DG:5:VAL:HG13	2.39	0.52
43:DI:109:ILE:HD12	43:DI:109:ILE:O	2.09	0.52
43:DI:94:ALA:HA	43:DI:97:ILE:HB	1.90	0.52
44:DN:13:TRP:O	44:DN:135:PRO:HD2	2.09	0.52
46:DP:19:VAL:HG23	46:DP:19:VAL:O	2.10	0.52
50:DT:51:ARG:HH11	50:DT:51:ARG:HG3	1.74	0.52
52:DV:78:LYS:HD3	52:DV:79:VAL:CA	2.40	0.52
53:DW:20:VAL:HG21	53:DW:47:VAL:HG21	1.92	0.52
1:AA:1372:U:H5''	9:AI:71:SER:CB	2.35	0.52
1:AA:286:G:H2'	1:AA:287:U:C6	2.44	0.52
1:AA:376:G:H2'	1:AA:377:G:C8	2.45	0.52
1:AA:393:A:O2'	1:AA:394:G:H5'	2.10	0.52
1:AA:402:G:H4'	1:AA:620:C:N3	2.24	0.52
1:AA:425:G:H2'	1:AA:426:G:C8	2.43	0.52
1:AA:527:G:H2'	1:AA:528:C:H5'	1.91	0.52
1:AA:643:C:H5'	8:AH:31:PHE:CE1	2.44	0.52
1:AA:814:A:H2'	1:AA:816:A:H5''	1.91	0.52
1:AA:973:G:H1'	10:AJ:54:PHE:CE1	2.44	0.52
2:AB:124:SER:OG	2:AB:125:PRO:HD2	2.08	0.52
4:AD:138:TYR:HD2	4:AD:139:ARG:N	2.07	0.52
4:AD:18:LYS:CE	4:AD:31:CYS:SG	2.98	0.52
5:AE:50:GLU:OE2	5:AE:51:VAL:HG23	2.10	0.52
8:AH:28:ALA:HA	8:AH:59:LEU:HG	1.92	0.52
1:AA:539:A:OP1	12:AL:114:LYS:HE2	2.09	0.52
12:AL:60:LEU:HD22	12:AL:60:LEU:N	2.24	0.52
1:AA:280:C:C4	17:AQ:91:ARG:NH2	2.78	0.52
19:AS:63:THR:HG22	19:AS:66:MET:HE2	1.91	0.52
21:AU:21:TYR:CD1	21:AU:21:TYR:N	2.75	0.52
23:AW:24:C:H2'	23:AW:25:U:O4'	2.09	0.52
26:B0:72:ARG:HB3	26:B0:75:LEU:HB3	1.92	0.52
27:B1:33:LYS:HG2	27:B1:34:THR:N	2.21	0.52
27:B1:64:ALA:O	27:B1:67:ILE:CG1	2.50	0.52
31:B5:15:ARG:HA	31:B5:18:ALA:CB	2.39	0.52
35:BA:1039:G:H3'	35:BA:1040:C:C6	2.44	0.52
35:BA:1431:U:O2'	35:BA:1432:C:H5'	2.10	0.52
35:BA:1516:C:H2'	35:BA:1517:G:H8	1.73	0.52
35:BA:1628:G:O2'	35:BA:1629:U:H5'	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1707:G:H1'	35:BA:1756:G:N3	2.24	0.52
35:BA:2250:G:C6	47:BQ:82:ARG:HD2	2.44	0.52
35:BA:2619:C:O2'	35:BA:2620:C:H5'	2.09	0.52
35:BA:1999:C:H4'	35:BA:2723:C:O2	2.08	0.52
35:BA:2807:G:C3'	35:BA:2808:U:H5''	2.40	0.52
35:BA:2828:C:O2'	35:BA:2829:C:H5'	2.09	0.52
35:BA:435:C:H2'	35:BA:436:C:H5'	1.92	0.52
35:BA:79:G:O2'	35:BA:80:G:H5'	2.09	0.52
37:BC:59:ARG:HH21	37:BC:199:HIS:CB	2.22	0.52
37:BC:214:VAL:C	37:BC:216:THR:N	2.62	0.52
37:BC:22:ILE:CG2	37:BC:25:ALA:HB2	2.40	0.52
38:BD:73:VAL:HG13	38:BD:120:GLY:CA	2.39	0.52
38:BD:127:VAL:HA	38:BD:193:VAL:HG13	1.91	0.52
38:BD:160:GLY:N	38:BD:196:VAL:HB	2.24	0.52
38:BD:39:LYS:HB2	38:BD:62:TYR:HB2	1.90	0.52
39:BE:128:SER:O	39:BE:129:HIS:CB	2.55	0.52
41:BG:98:ARG:O	41:BG:101:ILE:HG23	2.10	0.52
41:BG:138:GLN:HB3	41:BG:153:ARG:O	2.10	0.52
42:BH:138:LYS:HA	42:BH:141:VAL:HB	1.92	0.52
45:BO:16:ALA:CB	45:BO:43:VAL:HG13	2.36	0.52
46:BP:71:VAL:CB	46:BP:72:PRO:HD3	2.38	0.52
51:BU:64:ARG:NH2	51:BU:64:ARG:HA	2.25	0.52
52:BV:34:GLU:C	52:BV:62:LEU:HD12	2.29	0.52
53:BW:70:TYR:O	53:BW:107:LEU:HG	2.09	0.52
1:CA:1208:C:H2'	1:CA:1209:C:H6	1.73	0.52
1:CA:1321:C:H5''	1:CA:1322:C:H5''	1.89	0.52
1:CA:297:G:H2'	1:CA:299:G:OP2	2.09	0.52
3:CC:109:PRO:O	3:CC:115:LEU:HD12	2.08	0.52
4:CD:119:GLN:HE21	4:CD:123:HIS:CD2	2.27	0.52
4:CD:65:ARG:NH1	4:CD:72:GLU:N	2.57	0.52
5:CE:31:LEU:CD1	5:CE:129:ILE:HA	2.39	0.52
7:CG:20:ASP:HB3	7:CG:23:VAL:HG23	1.91	0.52
11:CK:69:ALA:O	11:CK:70:LYS:C	2.48	0.52
15:CO:9:GLN:O	15:CO:10:LYS:C	2.48	0.52
16:CP:14:ASN:H	16:CP:15:PRO:HD3	1.73	0.52
16:CP:1:MET:SD	16:CP:3:LYS:HE3	2.50	0.52
16:CP:49:LEU:CD1	16:CP:51:VAL:HG23	2.35	0.52
17:CQ:31:LEU:HG	17:CQ:32:TYR:CE2	2.44	0.52
17:CQ:86:GLU:O	17:CQ:88:TYR:N	2.42	0.52
18:CR:85:LEU:HG	18:CR:86:VAL:N	2.16	0.52
28:D2:22:GLU:C	28:D2:24:LEU:N	2.63	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:111:A:O2'	35:DA:112:U:H5'	2.08	0.52
35:DA:1132:A:H2'	35:DA:1133:U:C6	2.44	0.52
35:DA:1720:U:H2'	35:DA:1721:G:O4'	2.10	0.52
35:DA:172:C:O2	35:DA:172:C:H2'	2.09	0.52
35:DA:2040:C:H2'	35:DA:2041:U:H6	1.74	0.52
35:DA:2064:C:H1'	35:DA:2450:A:C6	2.45	0.52
35:DA:2465:C:O2'	35:DA:2466:C:H5'	2.10	0.52
35:DA:271(Q):G:H2'	35:DA:271(R):G:H8	1.74	0.52
35:DA:2805:G:H22	35:DA:2893:G:H1	1.57	0.52
35:DA:301:G:C4	35:DA:302:C:C5	2.97	0.52
35:DA:959:A:H2'	35:DA:960:A:C8	2.45	0.52
36:DB:20:C:H2'	36:DB:21:G:C5'	2.38	0.52
38:DD:248:SER:HB2	38:DD:249:PRO:HD2	1.92	0.52
39:DE:116:VAL:CG2	39:DE:122:PHE:CG	2.92	0.52
39:DE:52:LEU:HB2	39:DE:76:ARG:CB	2.34	0.52
39:DE:59:VAL:O	39:DE:62:PRO:HD2	2.09	0.52
41:DG:156:ASP:O	41:DG:157:ILE:O	2.26	0.52
45:DO:98:VAL:HG22	45:DO:99:PHE:N	2.25	0.52
47:DQ:119:ARG:HG2	47:DQ:120:ILE:CD1	2.36	0.52
49:DS:24:LEU:O	49:DS:86:ALA:HB3	2.10	0.52
50:DT:25:GLY:HA2	50:DT:92:GLY:HA3	1.91	0.52
50:DT:54:ARG:HA	50:DT:59:THR:OG1	2.09	0.52
1:AA:1229:A:H2'	1:AA:1230:C:H6	1.75	0.52
1:AA:189(F):U:O4	17:AQ:62:SER:HB3	2.09	0.52
1:AA:261:U:H3'	20:AT:79:ARG:HH12	1.75	0.52
1:AA:627:G:O2'	1:AA:628:G:H5'	2.09	0.52
2:AB:212:GLN:HE22	2:AB:216:SER:CB	2.22	0.52
3:AC:125:GLU:CG	3:AC:189:ALA:HA	2.38	0.52
7:AG:126:ASP:HB3	7:AG:132:GLY:CA	2.40	0.52
7:AG:140:ASP:OD1	7:AG:143:ARG:NH2	2.40	0.52
7:AG:15:ASP:OD2	7:AG:16:LEU:N	2.43	0.52
10:AJ:48:THR:HG23	10:AJ:62:HIS:N	2.24	0.52
10:AJ:85:LEU:O	10:AJ:87:THR:N	2.42	0.52
15:AO:82:ILE:HD13	15:AO:82:ILE:C	2.30	0.52
16:AP:49:LEU:HD22	16:AP:73:LEU:HD13	1.90	0.52
19:AS:36:ARG:HH22	19:AS:75:ALA:CB	2.12	0.52
29:B3:35:ARG:HG2	29:B3:37:LEU:HD21	1.90	0.52
34:B8:49:VAL:HG12	34:B8:53:PRO:HD3	1.91	0.52
35:BA:818:G:OP2	35:BA:1187:G:O6	2.27	0.52
33:B7:9:ARG:NH1	35:BA:1310:G:OP2	2.43	0.52
35:BA:172:C:O2	35:BA:172:C:H2'	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1791:A:N6	35:BA:1828:G:O2'	2.43	0.52
35:BA:1922:G:O2'	35:BA:1923:U:H5'	2.10	0.52
35:BA:2078:C:O2'	35:BA:2079:U:H5'	2.09	0.52
26:B0:16:SER:HB3	35:BA:2262:U:OP2	2.09	0.52
35:BA:251:A:H2'	35:BA:252:G:O4'	2.09	0.52
35:BA:2643:G:O2'	35:BA:2644:G:H5'	2.08	0.52
35:BA:2676:C:O2'	35:BA:2677:G:H5'	2.10	0.52
35:BA:325:G:H2'	35:BA:326:G:H8	1.74	0.52
35:BA:419:C:O2'	35:BA:420:C:H5'	2.10	0.52
35:BA:921:G:H2'	35:BA:922:U:H6	1.74	0.52
35:BA:848:G:N3	35:BA:933:A:H1'	2.24	0.52
36:BB:15:A:C3'	36:BB:16:G:H5'	2.40	0.52
38:BD:264:LYS:HE3	38:BD:266:SER:HB2	1.91	0.52
39:BE:2:LYS:CD	39:BE:95:ILE:HG22	2.39	0.52
40:BF:41:LEU:HA	40:BF:44:ARG:HD3	1.92	0.52
41:BG:112:PRO:O	41:BG:113:ARG:HA	2.09	0.52
42:BH:105:LEU:CD2	42:BH:113:VAL:HB	2.38	0.52
44:BN:62:VAL:CG2	44:BN:66:LYS:HB2	2.40	0.52
44:BN:72:TYR:N	44:BN:85:ILE:O	2.42	0.52
45:BO:112:MET:O	45:BO:113:LYS:C	2.48	0.52
46:BP:122:PRO:HA	46:BP:141:ALA:O	2.10	0.52
49:BS:72:ALA:O	49:BS:73:LEU:C	2.47	0.52
35:BA:17:G:HO2'	51:BU:25:TRP:HZ3	1.55	0.52
52:BV:14:VAL:HG11	52:BV:98:GLU:HG3	1.91	0.52
35:BA:58:G:OP1	54:BX:72:LYS:CB	2.56	0.52
56:BZ:5:LEU:N	56:BZ:59:LEU:HD23	2.25	0.52
1:CA:1365:G:C2'	1:CA:1366:C:H5'	2.39	0.52
1:CA:277:C:O2'	1:CA:278:G:H5'	2.10	0.52
1:CA:324:G:N2	1:CA:326:G:H3'	2.24	0.52
1:CA:406:G:C2	1:CA:407:G:C8	2.97	0.52
1:CA:601:C:H2'	1:CA:602:A:C8	2.42	0.52
1:CA:727:G:N1	1:CA:731:G:C6	2.77	0.52
1:CA:764:C:H2'	1:CA:765:G:O4'	2.08	0.52
1:CA:835:U:OP1	18:CR:61:LYS:HB2	2.10	0.52
1:CA:853:G:H2'	1:CA:854:G:H8	1.73	0.52
1:CA:960:U:H2'	1:CA:960:U:O2	2.09	0.52
1:CA:998:G:H2'	1:CA:999:C:C6	2.43	0.52
2:CB:167:PRO:O	2:CB:168:THR:C	2.48	0.52
3:CC:5:ILE:O	3:CC:5:ILE:HD13	2.09	0.52
4:CD:172:PRO:HD2	4:CD:173:TRP:CZ3	2.45	0.52
4:CD:61:LYS:HG2	4:CD:75:PHE:HE2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:101:ILE:CD1	5:CE:118:ILE:O	2.57	0.52
5:CE:15:ARG:O	5:CE:15:ARG:HG2	2.10	0.52
5:CE:80:ILE:HG13	5:CE:91:LEU:HB2	1.92	0.52
8:CH:21:LYS:CG	8:CH:22:GLU:N	2.73	0.52
8:CH:7:ALA:O	8:CH:11:THR:N	2.38	0.52
10:CJ:34:VAL:HG12	10:CJ:35:SER:N	2.24	0.52
11:CK:69:ALA:O	11:CK:73:MET:N	2.31	0.52
13:CM:67:GLU:O	13:CM:69:GLU:N	2.42	0.52
15:CO:70:LEU:HG	15:CO:78:TYR:HB2	1.90	0.52
17:CQ:47:PRO:HG2	17:CQ:48:GLU:OE1	2.08	0.52
17:CQ:85:VAL:O	17:CQ:89:LEU:HB2	2.10	0.52
18:CR:75:ILE:O	18:CR:75:ILE:HG22	2.08	0.52
19:CS:16:LEU:N	19:CS:16:LEU:HD12	2.25	0.52
20:CT:25:ARG:HG3	20:CT:25:ARG:HH11	1.75	0.52
20:CT:63:ILE:O	20:CT:66:ALA:N	2.43	0.52
28:D2:20:GLU:C	28:D2:22:GLU:N	2.61	0.52
35:DA:81:G:H1	35:DA:105:C:H42	1.55	0.52
35:DA:1644:C:HO2'	35:DA:1645:G:H5'	1.73	0.52
35:DA:2522:U:H2'	35:DA:2523:G:H5''	1.91	0.52
35:DA:2649:U:O2'	35:DA:2650:U:H5'	2.09	0.52
35:DA:2684:U:H2'	35:DA:2685:G:O4'	2.09	0.52
35:DA:2728:U:H2'	35:DA:2728:U:O2	2.09	0.52
35:DA:2790:A:N3	35:DA:2790:A:H2'	2.24	0.52
35:DA:729:G:N7	38:DD:208:LYS:HB2	2.24	0.52
36:DB:31:C:H42	36:DB:51:G:H1	1.56	0.52
36:DB:90:A:C8	36:DB:91:C:H1'	2.44	0.52
35:DA:1902:C:O2'	38:DD:244:ARG:HB2	2.09	0.52
39:DE:68:ALA:C	39:DE:70:ALA:H	2.12	0.52
40:DF:153:SER:HA	40:DF:172:TRP:O	2.09	0.52
40:DF:45:ARG:HG3	40:DF:46:ARG:H	1.75	0.52
44:DN:42:TRP:HA	44:DN:42:TRP:CE3	2.44	0.52
49:DS:36:TYR:HA	49:DS:52:SER:CB	2.39	0.52
50:DT:102:ILE:HB	50:DT:110:ILE:HD11	1.91	0.52
45:DO:80:ASP:HB2	50:DT:71:GLY:O	2.10	0.52
50:DT:94:ALA:CB	50:DT:99:LEU:HD23	2.40	0.52
54:DX:53:LYS:NZ	54:DX:55:ASN:ND2	2.58	0.52
54:DX:54:VAL:HG13	54:DX:78:LYS:O	2.09	0.52
56:DZ:168:GLU:O	56:DZ:169:GLU:O	2.28	0.52
1:AA:1221:G:OP1	19:AS:36:ARG:HD3	2.08	0.52
1:AA:1365:G:C2'	1:AA:1366:C:H5'	2.40	0.52
1:AA:1521:G:H2'	1:AA:1522:U:H6	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:197:A:N3	1:AA:198:G:H1'	2.25	0.52
1:AA:337:C:O2'	1:AA:338:A:H5'	2.10	0.52
1:AA:423:G:H2'	1:AA:424:G:O4'	2.09	0.52
1:AA:59:A:N3	1:AA:59:A:H2'	2.24	0.52
1:AA:640:A:O2'	1:AA:641:U:H5'	2.10	0.52
1:AA:569:C:N4	1:AA:881:G:H1	2.07	0.52
1:AA:973:G:C4	10:AJ:55:LYS:NZ	2.68	0.52
2:AB:221:LEU:O	2:AB:221:LEU:HD13	2.09	0.52
3:AC:106:VAL:HG12	3:AC:108:ASN:H	1.74	0.52
3:AC:147:LYS:HB3	3:AC:203:PHE:CE2	2.45	0.52
4:AD:58:LEU:O	4:AD:60:GLU:N	2.43	0.52
4:AD:61:LYS:HA	4:AD:203:VAL:CG2	2.32	0.52
1:AA:875:C:O2'	8:AH:14:ARG:NH1	2.43	0.52
12:AL:46:LYS:CG	12:AL:47:LYS:H	2.08	0.52
12:AL:7:ILE:O	12:AL:10:LEU:HB2	2.10	0.52
12:AL:86:ARG:CG	12:AL:87:GLY:N	2.73	0.52
15:AO:12:ILE:C	15:AO:14:GLU:H	2.13	0.52
19:AS:22:LEU:C	19:AS:24:ALA:H	2.11	0.52
20:AT:26:ASN:ND2	20:AT:26:ASN:N	2.56	0.52
27:B1:23:LYS:O	27:B1:37:ILE:HD12	2.09	0.52
27:B1:87:PRO:CG	27:B1:88:LYS:N	2.73	0.52
35:BA:1186:G:C2'	35:BA:1187:G:H5'	2.39	0.52
35:BA:1279:G:H2'	35:BA:1280:G:H8	1.74	0.52
35:BA:1301:A:H2	35:BA:1626:G:H21	1.57	0.52
35:BA:1300:U:O2	35:BA:1626:G:C4	2.63	0.52
35:BA:1665:A:H1'	45:BO:1:MET:HE3	1.90	0.52
35:BA:1720:U:H2'	35:BA:1721:G:H5'	1.91	0.52
35:BA:1776:G:C2	35:BA:1777:U:C6	2.96	0.52
35:BA:1814:G:H2'	35:BA:1815:A:N7	2.24	0.52
35:BA:1827:C:O2'	35:BA:1828:G:H5'	2.10	0.52
35:BA:2069:G:C2'	35:BA:2070:G:H5'	2.39	0.52
35:BA:2181:G:H2'	35:BA:2182:G:C8	2.45	0.52
35:BA:2250:G:C8	35:BA:2496:C:H5''	2.44	0.52
35:BA:2593:U:H2'	35:BA:2594:C:C6	2.44	0.52
35:BA:268:C:C2'	35:BA:268:C:O2	2.55	0.52
35:BA:554:U:C2'	35:BA:555:U:H5'	2.40	0.52
35:BA:813:U:H2'	35:BA:814:C:H6	1.72	0.52
35:BA:840:C:O2'	35:BA:841:A:H5'	2.09	0.52
35:BA:869:G:O2'	35:BA:870:A:H5'	2.09	0.52
35:BA:955:C:H5'	35:BA:956:G:OP2	2.09	0.52
38:BD:142:VAL:HG22	38:BD:143:HIS:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:56:PRO:O	39:BE:57:LYS:O	2.27	0.52
40:BF:123:LEU:HD12	40:BF:124:LEU:H	1.74	0.52
40:BF:9:ILE:O	40:BF:128:ALA:HB2	2.10	0.52
41:BG:167:GLU:H	41:BG:167:GLU:CD	2.13	0.52
42:BH:149:ARG:CA	42:BH:162:ILE:HD11	2.38	0.52
43:BI:68:LEU:CD2	43:BI:136:VAL:HG11	2.40	0.52
44:BN:42:TRP:HE3	44:BN:42:TRP:HA	1.73	0.52
46:BP:58:THR:O	46:BP:61:ARG:NE	2.43	0.52
48:BR:117:VAL:O	48:BR:118:GLU:HB2	2.10	0.52
48:BR:52:ILE:CD1	48:BR:79:LEU:HD21	2.40	0.52
55:BY:98:VAL:O	55:BY:98:VAL:HG12	2.10	0.52
1:CA:1037:C:H2'	1:CA:1038:C:C6	2.45	0.52
1:CA:1155:G:O2'	1:CA:1156:G:H5'	2.09	0.52
1:CA:185:A:H61	1:CA:192:U:H3	1.58	0.52
1:CA:355:C:C4	1:CA:356:A:N7	2.77	0.52
1:CA:486:U:O2'	1:CA:487:A:H5'	2.09	0.52
1:CA:501:C:H2'	1:CA:502:G:C8	2.44	0.52
1:CA:503:C:O2'	1:CA:504:C:H5'	2.10	0.52
1:CA:741:G:H5'	15:CO:39:LEU:HD21	1.91	0.52
1:CA:865:A:C2	1:CA:918:A:H4'	2.44	0.52
1:CA:569:C:N4	1:CA:881:G:H1	2.07	0.52
2:CB:136:VAL:O	2:CB:140:HIS:HB2	2.09	0.52
2:CB:69:LEU:HB2	2:CB:159:PRO:CG	2.39	0.52
4:CD:65:ARG:HB2	4:CD:75:PHE:CE2	2.44	0.52
4:CD:92:VAL:O	4:CD:95:GLY:N	2.43	0.52
5:CE:92:LYS:HB2	5:CE:119:LEU:HB2	1.91	0.52
8:CH:129:VAL:HG23	8:CH:130:GLY:N	2.15	0.52
12:CL:38:THR:HG23	12:CL:39:VAL:N	2.24	0.52
14:CN:53:LEU:HB3	14:CN:56:VAL:HG21	1.90	0.52
15:CO:66:LEU:HD12	15:CO:66:LEU:H	1.73	0.52
17:CQ:71:PHE:N	17:CQ:71:PHE:HD2	2.07	0.52
19:CS:33:THR:HG21	19:CS:51:VAL:HG22	1.92	0.52
20:CT:81:LYS:O	20:CT:85:MET:HG2	2.10	0.52
25:CY:63:PRO:O	25:CY:65:THR:N	2.36	0.52
28:D2:48:HIS:O	28:D2:52:ASP:HB3	2.09	0.52
33:D7:29:LYS:O	33:D7:33:ARG:N	2.39	0.52
35:DA:1424:G:O2'	35:DA:1425:G:H5'	2.09	0.52
35:DA:149:A:C2	35:DA:150:C:C2	2.97	0.52
35:DA:1668:A:N3	35:DA:1670:C:C4	2.78	0.52
35:DA:1917:U:H2'	35:DA:1918:A:C8	2.44	0.52
35:DA:2066:C:O2'	35:DA:2067:G:H5'	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:D1:37:ILE:HD12	35:DA:2080:G:P	2.50	0.52
35:DA:2713:A:C3'	35:DA:2714:G:H5'	2.38	0.52
35:DA:407:G:H2'	35:DA:408:G:C8	2.45	0.52
35:DA:745:G:H2'	35:DA:746:A:H5'	1.90	0.52
35:DA:785:G:C4	35:DA:786:C:C5	2.97	0.52
35:DA:811:U:H1'	35:DA:1251:C:H5''	1.91	0.52
37:DC:23:ASP:C	37:DC:25:ALA:H	2.13	0.52
37:DC:77:ILE:HG12	37:DC:77:ILE:O	2.08	0.52
38:DD:44:ASN:ND2	38:DD:47:GLY:O	2.42	0.52
39:DE:5:LEU:CD2	39:DE:197:ILE:HG22	2.40	0.52
40:DF:155:LEU:CD1	40:DF:174:VAL:HB	2.40	0.52
41:DG:106:LEU:C	41:DG:108:ASN:N	2.63	0.52
41:DG:132:ASN:ND2	41:DG:157:ILE:O	2.42	0.52
43:DI:98:ALA:O	43:DI:109:ILE:HG21	2.10	0.52
43:DI:49:ALA:HA	43:DI:52:ARG:CG	2.37	0.52
35:DA:196:A:C5'	46:DP:46:LYS:HZ1	2.22	0.52
35:DA:2839:G:H21	48:DR:92:GLY:HA3	1.74	0.52
50:DT:23:ARG:C	50:DT:25:GLY:H	2.05	0.52
1:CA:346:G:H5''	50:DT:35:LYS:HZ1	1.75	0.52
52:DV:23:GLU:O	52:DV:24:LYS:O	2.27	0.52
54:DX:62:LYS:HD2	54:DX:68:ARG:HD2	1.92	0.52
54:DX:76:ARG:C	54:DX:76:ARG:HD3	2.30	0.52
54:DX:82:GLN:OE1	54:DX:83:VAL:N	2.41	0.52
55:DY:86:ARG:NH2	55:DY:95:LYS:HZ3	2.08	0.52
56:DZ:144:LEU:HD12	56:DZ:149:SER:HA	1.91	0.52
1:AA:1427:U:O2'	1:AA:1428:A:H5'	2.10	0.52
1:AA:885:G:H1	1:AA:912:C:H42	1.57	0.52
1:AA:914:A:O2'	1:AA:915:A:H5'	2.10	0.52
2:AB:114:ARG:HA	2:AB:117:GLU:OE1	2.10	0.52
6:AF:100:ASN:HB3	18:AR:28:GLU:HA	1.92	0.52
8:AH:53:VAL:O	8:AH:54:ASP:HB2	2.09	0.52
9:AI:4:TYR:CB	9:AI:19:LEU:HB2	2.37	0.52
10:AJ:7:LYS:O	10:AJ:96:ILE:HA	2.09	0.52
13:AM:64:TRP:HE1	13:AM:66:LEU:HD12	1.75	0.52
16:AP:4:ILE:HB	16:AP:66:PRO:CB	2.32	0.52
17:AQ:29:HIS:HA	17:AQ:36:ILE:HD11	1.91	0.52
25:AY:126:ARG:O	25:AY:129:ILE:HB	2.10	0.52
26:B0:19:LYS:C	26:B0:20:ARG:HD3	2.30	0.52
26:B0:53:MET:HG3	26:B0:59:LEU:HD23	1.91	0.52
27:B1:73:LEU:HD21	27:B1:94:LEU:CD2	2.36	0.52
29:B3:32:GLN:HB2	35:BA:1158:C:H4'	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:B3:54:VAL:HG12	29:B3:55:ARG:N	2.24	0.52
34:B8:29:LYS:HZ3	34:B8:44:LYS:HB2	1.74	0.52
35:BA:1036:G:O2'	35:BA:1037:G:H5'	2.09	0.52
35:BA:1151:G:H5''	51:BU:81:HIS:CE1	2.45	0.52
35:BA:991:C:N4	35:BA:1163:G:H1	2.07	0.52
35:BA:1181:C:H2'	35:BA:1182:A:C8	2.45	0.52
35:BA:1445(A):C:H2'	35:BA:1446:C:H6	1.74	0.52
35:BA:1720:U:H2'	35:BA:1721:G:O4'	2.09	0.52
35:BA:2298:A:H62	35:BA:2318:G:H8	1.54	0.52
35:BA:2313:C:H5'	35:BA:2313:C:C6	2.44	0.52
35:BA:2413:G:H21	46:BP:70:GLN:NE2	2.07	0.52
35:BA:265:A:H1'	35:BA:266:G:C1'	2.38	0.52
35:BA:445:C:H5''	51:BU:3:ARG:HB3	1.90	0.52
35:BA:738:G:C6	35:BA:739:G:C2	2.98	0.52
35:BA:796:C:O2'	35:BA:797:C:H5'	2.09	0.52
38:BD:9:TYR:HD2	38:BD:10:THR:HG22	1.73	0.52
40:BF:192:LEU:HD21	40:BF:194:MET:HE2	1.91	0.52
40:BF:2:LYS:CG	40:BF:25:PRO:HB2	2.22	0.52
41:BG:139:LEU:HD22	41:BG:146:TYR:CE1	2.44	0.52
41:BG:170:ARG:HH22	41:BG:182:LYS:NZ	2.07	0.52
43:BI:9:LEU:H	43:BI:13:GLY:HA2	1.74	0.52
43:BI:94:ALA:HA	43:BI:97:ILE:HB	1.91	0.52
45:BO:26:LYS:HB2	45:BO:30:ALA:CB	2.39	0.52
45:BO:85:VAL:HG12	45:BO:86:ILE:N	2.25	0.52
46:BP:138:LEU:HD23	46:BP:142:GLY:HA3	1.92	0.52
47:BQ:35:VAL:HG22	47:BQ:100:GLY:O	2.09	0.52
50:BT:113:LYS:C	50:BT:114:LEU:HD23	2.30	0.52
56:BZ:112:ARG:O	56:BZ:113:ALA:HB2	2.09	0.52
56:BZ:4:ARG:HB3	56:BZ:60:GLU:HG3	1.92	0.52
1:CA:1062:U:H2'	1:CA:1063:C:C5	2.44	0.52
1:CA:1332:A:H2'	1:CA:1333:A:C8	2.45	0.52
1:CA:1333:A:C2	1:CA:1334:G:H1'	2.45	0.52
1:CA:1354:C:O2'	1:CA:1355:G:H5'	2.10	0.52
1:CA:184:G:O2'	1:CA:185:A:H5'	2.09	0.52
1:CA:401:C:H6	1:CA:401:C:O5'	1.93	0.52
1:CA:509:A:H2'	1:CA:510:A:C8	2.45	0.52
1:CA:814:A:H2'	1:CA:816:A:H5''	1.92	0.52
1:CA:950:U:H1'	1:CA:971:G:N7	2.24	0.52
2:CB:212:GLN:HE22	2:CB:216:SER:HB2	1.75	0.52
4:CD:170:VAL:HG22	4:CD:171:GLY:H	1.75	0.52
4:CD:170:VAL:HG13	4:CD:174:LEU:HB2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:62:GLN:O	4:CD:63:LYS:C	2.48	0.52
17:CQ:43:LEU:HD12	17:CQ:68:ARG:HG3	1.92	0.52
19:CS:70:LYS:H	19:CS:70:LYS:HD2	1.74	0.52
20:CT:58:LYS:O	20:CT:62:LEU:HB2	2.09	0.52
23:CW:64:G:H2'	23:CW:65:G:O4'	2.10	0.52
27:D1:46:LEU:HA	35:DA:396:G:O3'	2.10	0.52
28:D2:27:GLU:C	28:D2:29:LYS:N	2.56	0.52
28:D2:57:ILE:HD11	28:D2:59:ARG:HH11	1.73	0.52
35:DA:1141:U:H5''	35:DA:1142(A):A:O4'	2.10	0.52
35:DA:1150:C:C2'	35:DA:1151:G:H5'	2.39	0.52
35:DA:154:G:H2'	35:DA:154(A):C:O2	2.08	0.52
35:DA:1612:C:H2'	35:DA:1613:G:O5'	2.09	0.52
35:DA:1791:A:N6	35:DA:1828:G:O2'	2.42	0.52
35:DA:1860:G:H5''	37:DC:204:ALA:O	2.10	0.52
35:DA:1986:A:C2'	35:DA:1987:G:H5''	2.40	0.52
35:DA:2383:G:O2'	35:DA:2384:G:H5'	2.10	0.52
35:DA:2809:A:C2	35:DA:2892:A:N3	2.77	0.52
38:DD:130:ALA:HB2	38:DD:192:THR:HB	1.91	0.52
39:DE:142:GLY:C	39:DE:143:ASN:HD22	2.13	0.52
39:DE:152:LYS:HZ3	44:DN:78:TYR:CB	2.22	0.52
39:DE:201:THR:CG2	39:DE:202:LYS:N	2.72	0.52
39:DE:47:VAL:CG2	39:DE:84:PHE:O	2.57	0.52
40:DF:123:LEU:HD12	40:DF:124:LEU:H	1.75	0.52
40:DF:34:TRP:HB3	46:DP:11:GLY:HA3	1.91	0.52
40:DF:89:VAL:HG12	40:DF:90:PHE:H	1.75	0.52
35:DA:2304:G:O2'	41:DG:133:LEU:HA	2.10	0.52
41:DG:132:ASN:HD22	41:DG:133:LEU:N	2.08	0.52
41:DG:35:GLU:OE2	41:DG:160:VAL:HG11	2.09	0.52
42:DH:102:ALA:HB2	42:DH:117:PRO:CD	2.21	0.52
42:DH:130:ARG:HB2	42:DH:130:ARG:NH1	2.25	0.52
44:DN:42:TRP:HA	44:DN:42:TRP:HE3	1.74	0.52
45:DO:26:LYS:HB2	45:DO:30:ALA:CB	2.40	0.52
45:DO:88:ASN:O	45:DO:91:LEU:N	2.42	0.52
49:DS:32:LEU:O	49:DS:33:LYS:HB2	2.09	0.52
45:DO:104:ARG:CZ	50:DT:33:LYS:HD2	2.39	0.52
54:DX:49:VAL:CG1	54:DX:50:LYS:N	2.70	0.52
56:DZ:102:LEU:HB2	56:DZ:122:ARG:O	2.10	0.52
1:AA:240:C:H2'	1:AA:241:C:H6	1.74	0.52
1:AA:261:U:C6	20:AT:79:ARG:NH1	2.78	0.52
1:AA:437:U:C2'	1:AA:438:G:H5'	2.40	0.52
2:AB:167:PRO:HD2	2:AB:188:ALA:HB2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:182:ILE:HG23	3:AC:203:PHE:N	2.25	0.52
6:AF:14:LEU:HD13	6:AF:15:ASP:O	2.10	0.52
7:AG:21:VAL:HG23	7:AG:22:LEU:N	2.24	0.52
11:AK:125:PHE:N	11:AK:125:PHE:CD1	2.78	0.52
12:AL:70:ILE:N	12:AL:70:ILE:CD1	2.66	0.52
16:AP:22:THR:CG2	16:AP:32:TYR:HA	2.40	0.52
18:AR:19:LYS:O	18:AR:20:ALA:CB	2.58	0.52
19:AS:27:GLU:O	19:AS:28:LYS:HD2	2.09	0.52
19:AS:70:LYS:H	19:AS:70:LYS:HD2	1.74	0.52
20:AT:26:ASN:CB	20:AT:71:THR:HG23	2.36	0.52
22:AV:27:G:H2'	22:AV:28:G:H8	1.74	0.52
25:AY:133:ARG:HD3	25:AY:165:THR:CG2	2.28	0.52
27:B1:31:GLY:O	35:BA:2397:G:H5'	2.09	0.52
27:B1:19:GLN:O	27:B1:42:GLN:HB3	2.09	0.52
28:B2:12:GLU:O	28:B2:12:GLU:CD	2.48	0.52
28:B2:16:LEU:O	28:B2:20:GLU:HG3	2.10	0.52
28:B2:51:ARG:CD	28:B2:51:ARG:O	2.58	0.52
34:B8:49:VAL:CB	34:B8:53:PRO:HD3	2.39	0.52
34:B8:56:GLU:C	34:B8:58:ILE:N	2.62	0.52
35:BA:1014:U:H2'	35:BA:1015:G:C8	2.45	0.52
35:BA:1980:G:C2	35:BA:1982:C:C4	2.97	0.52
35:BA:2115:G:O2'	35:BA:2116:G:H5''	2.10	0.52
35:BA:2119:A:H3'	35:BA:2120:G:H5''	1.91	0.52
35:BA:2260:C:H2'	35:BA:2261:C:H6	1.74	0.52
35:BA:2290:G:H2'	35:BA:2291:U:O4'	2.10	0.52
35:BA:2368:C:H2'	35:BA:2369:A:C8	2.44	0.52
35:BA:236:C:O2'	35:BA:237:C:H5'	2.09	0.52
35:BA:2259:G:H1'	35:BA:2427:C:O2	2.10	0.52
35:BA:2536:G:C6	35:BA:2537:U:C4	2.97	0.52
35:BA:271(F):C:H2'	35:BA:271(G):C:H6	1.75	0.52
35:BA:271(L):U:H4'	35:BA:271(M):G:N7	2.24	0.52
35:BA:474:G:H4'	35:BA:475:U:OP1	2.10	0.52
35:BA:711:G:O2'	35:BA:712:G:H5'	2.10	0.52
35:BA:911:A:H5''	35:BA:912:C:H5''	1.90	0.52
36:BB:50:G:OP2	49:BS:62:LYS:HG3	2.09	0.52
38:BD:92:ILE:HA	38:BD:107:ALA:CB	2.40	0.52
38:BD:14:ARG:CG	38:BD:14:ARG:HH11	2.23	0.52
38:BD:166:GLN:N	38:BD:166:GLN:NE2	2.58	0.52
38:BD:54:ARG:H	38:BD:218:ARG:HG3	1.74	0.52
38:BD:65:ILE:HD11	38:BD:67:PHE:CE1	2.44	0.52
40:BF:125:LEU:HB3	40:BF:196:LEU:CD2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:114:ILE:HG22	41:BG:115:ARG:H	1.74	0.52
41:BG:133:LEU:HD13	41:BG:134:GLY:O	2.10	0.52
41:BG:29:TRP:C	41:BG:31:VAL:H	2.12	0.52
41:BG:83:ARG:O	41:BG:85:GLY:O	2.28	0.52
42:BH:137:ASP:O	42:BH:138:LYS:CB	2.57	0.52
43:BI:127:VAL:O	43:BI:127:VAL:HG12	2.09	0.52
44:BN:107:LEU:HD12	44:BN:108:PRO:C	2.30	0.52
46:BP:6:LEU:O	46:BP:6:LEU:HD23	2.10	0.52
47:BQ:35:VAL:HG22	47:BQ:101:ARG:O	2.09	0.52
49:BS:26:LEU:CD2	49:BS:26:LEU:O	2.57	0.52
49:BS:36:TYR:HA	49:BS:52:SER:CB	2.40	0.52
50:BT:100:TYR:C	50:BT:102:ILE:N	2.63	0.52
50:BT:3:ARG:O	50:BT:4:GLY:C	2.48	0.52
50:BT:51:ARG:HH11	50:BT:51:ARG:HG3	1.74	0.52
51:BU:88:ILE:HA	51:BU:90:VAL:HG23	1.92	0.52
51:BU:92:ARG:NH2	52:BV:10:LYS:HA	2.25	0.52
52:BV:3:ALA:O	52:BV:13:ARG:HA	2.10	0.52
56:BZ:108:PRO:CB	56:BZ:142:SER:O	2.58	0.52
1:CA:1207:G:O2'	1:CA:1208:C:H5'	2.10	0.52
1:CA:437:U:C2'	1:CA:438:G:H5'	2.40	0.52
1:CA:577:G:C4	1:CA:816:A:C2	2.98	0.52
1:CA:70:G:H2'	1:CA:71:C:C6	2.45	0.52
1:CA:939:G:H5''	7:CG:102:ARG:HH12	1.74	0.52
4:CD:15:GLU:C	4:CD:17:VAL:H	2.11	0.52
4:CD:60:GLU:O	4:CD:61:LYS:C	2.47	0.52
8:CH:40:ALA:HB2	8:CH:45:ILE:CD1	2.39	0.52
9:CI:50:LEU:O	9:CI:55:ALA:HB3	2.10	0.52
9:CI:114:TYR:HE1	10:CJ:60:ARG:H	1.56	0.52
7:CG:153:HIS:CE1	11:CK:57:THR:HG23	2.44	0.52
17:CQ:95:TYR:C	17:CQ:97:SER:N	2.63	0.52
19:CS:60:VAL:O	19:CS:60:VAL:HG13	2.09	0.52
25:CY:14:MET:HE3	25:CY:132:ILE:HB	1.92	0.52
25:CY:155:LYS:O	25:CY:158:GLU:HB3	2.10	0.52
29:D3:56:VAL:CG1	29:D3:57:GLU:H	2.05	0.52
34:D8:34:TRP:CZ3	34:D8:41:ILE:HG23	2.44	0.52
34:D8:56:GLU:O	34:D8:57:ARG:C	2.47	0.52
35:DA:1022:G:O2'	35:DA:1023:U:P	2.67	0.52
35:DA:991:C:N4	35:DA:1163:G:H1	2.07	0.52
35:DA:1658:C:H2'	35:DA:1659:U:C6	2.44	0.52
35:DA:1748:G:H8	35:DA:1748:G:H5'	1.73	0.52
35:DA:1899:G:O2'	35:DA:1900:A:H5''	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:237:C:O2'	35:DA:238:C:H5'	2.09	0.52
35:DA:2840:C:H2'	35:DA:2841:C:C6	2.44	0.52
35:DA:2845:G:H2'	35:DA:2846:G:H8	1.75	0.52
35:DA:354:G:O5'	35:DA:354:G:H8	1.93	0.52
35:DA:58:G:OP1	54:DX:72:LYS:CB	2.58	0.52
35:DA:782:A:C2	38:DD:226:MET:HE2	2.44	0.52
37:DC:44:HIS:HD2	37:DC:175:VAL:N	2.08	0.52
38:DD:24:ILE:O	38:DD:26:LYS:N	2.43	0.52
39:DE:103:ASP:CG	39:DE:201:THR:HA	2.29	0.52
39:DE:36:ARG:HA	39:DE:46:ALA:O	2.10	0.52
40:DF:125:LEU:HD11	40:DF:199:TRP:CG	2.45	0.52
40:DF:170:LEU:HD23	40:DF:173:VAL:CG2	2.39	0.52
41:DG:125:PHE:HE2	41:DG:173:LEU:CD1	2.23	0.52
35:DA:2312:U:OP1	41:DG:74:LYS:N	2.43	0.52
44:DN:107:LEU:HD12	44:DN:108:PRO:C	2.30	0.52
44:DN:58:ASP:C	44:DN:60:ILE:N	2.56	0.52
45:DO:1:MET:CG	45:DO:32:TYR:HD2	2.23	0.52
46:DP:107:LYS:O	46:DP:109:GLY:N	2.40	0.52
46:DP:40:SER:C	46:DP:41:ARG:HD2	2.30	0.52
49:DS:107:GLU:HG3	49:DS:108:GLY:N	2.25	0.52
53:DW:70:TYR:O	53:DW:107:LEU:HG	2.10	0.52
35:DA:1614:A:C6	53:DW:93:ALA:HB2	2.45	0.52
54:DX:29:TRP:HA	54:DX:29:TRP:HE3	1.75	0.52
54:DX:53:LYS:HZ2	54:DX:55:ASN:ND2	2.08	0.52
55:DY:49:VAL:O	55:DY:50:ARG:HB2	2.09	0.52
56:DZ:100:VAL:HG22	56:DZ:136:PHE:HA	1.91	0.52
56:DZ:61:LEU:O	56:DZ:63:ASP:N	2.39	0.52
1:AA:1253:G:O2'	1:AA:1254:C:H5'	2.10	0.52
1:AA:380:G:N1	1:AA:384:G:C6	2.78	0.52
1:AA:960:U:O2	1:AA:960:U:H2'	2.10	0.52
2:AB:9:GLU:O	2:AB:13:ALA:HB2	2.10	0.52
2:AB:74:LYS:CG	2:AB:77:ALA:HB3	2.39	0.52
4:AD:202:LEU:C	4:AD:204:ILE:N	2.62	0.52
4:AD:62:GLN:HB3	4:AD:66:ARG:CZ	2.40	0.52
7:AG:88:PRO:O	7:AG:89:MET:HB3	2.09	0.52
8:AH:122:ARG:NH1	8:AH:122:ARG:HB2	2.25	0.52
10:AJ:33:GLN:O	10:AJ:75:ILE:HG12	2.10	0.52
18:AR:22:VAL:HA	18:AR:25:THR:HG1	1.73	0.52
20:AT:30:LYS:HZ3	20:AT:72:LEU:HD21	1.75	0.52
23:AW:41:C:H2'	23:AW:42:C:C6	2.45	0.52
25:AY:80:GLU:C	25:AY:82:ALA:N	2.63	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:48:LYS:HD2	27:B1:61:ARG:HG2	1.91	0.52
35:BA:1142(A):A:C5	35:BA:1144:G:C5	2.98	0.52
35:BA:1341:U:OP1	35:BA:1397:U:N3	2.41	0.52
35:BA:139:G:H1	35:BA:142(A):C:H42	1.55	0.52
35:BA:1604:C:O2'	35:BA:1605:C:H5'	2.10	0.52
35:BA:1766:U:O2'	35:BA:1767:C:H5'	2.09	0.52
35:BA:1907:G:O2'	35:BA:1908:C:H5'	2.09	0.52
35:BA:1982:C:C2	35:BA:1983:C:C5	2.98	0.52
35:BA:199:A:N3	35:BA:2433:A:C2	2.77	0.52
35:BA:2464:C:O2'	35:BA:2465:C:H6	1.93	0.52
35:BA:2850:A:C2'	35:BA:2851:A:H8	2.16	0.52
35:BA:2889:C:O2	35:BA:2889:C:H2'	2.09	0.52
35:BA:815:C:OP2	52:BV:84:LYS:HE3	2.09	0.52
35:BA:85:G:O5'	55:BY:30:VAL:HB	2.08	0.52
35:BA:920:G:H2'	35:BA:921:G:C8	2.42	0.52
36:BB:104:U:H2'	36:BB:105:A:H8	1.74	0.52
37:BC:73:ARG:HA	37:BC:92:ASP:OD1	2.10	0.52
37:BC:76:ALA:H	37:BC:94:VAL:HG13	1.74	0.52
38:BD:211:ARG:O	38:BD:212:SER:C	2.48	0.52
35:BA:2632:A:N3	39:BE:61:ARG:NH1	2.57	0.52
39:BE:50:GLY:CA	39:BE:74:PRO:HG3	2.40	0.52
40:BF:171:PRO:HG2	40:BF:172:TRP:H	1.74	0.52
40:BF:89:VAL:C	40:BF:91:GLY:H	2.13	0.52
41:BG:125:PHE:CD2	41:BG:131:TYR:HB2	2.45	0.52
41:BG:171:ALA:O	41:BG:172:LEU:C	2.48	0.52
42:BH:45:VAL:O	42:BH:45:VAL:HG12	2.09	0.52
44:BN:97:ARG:HB3	44:BN:101:HIS:HD2	1.75	0.52
44:BN:26:LEU:HD21	44:BN:99:LEU:HD11	1.91	0.52
46:BP:47:ASP:HB2	46:BP:51:PHE:CD2	2.44	0.52
47:BQ:101:ARG:HG2	47:BQ:102:VAL:N	2.25	0.52
51:BU:24:TYR:HB2	51:BU:29:SER:CB	2.40	0.52
51:BU:74:LEU:HD11	51:BU:79:PHE:HB2	1.92	0.52
53:BW:29:LEU:HD23	53:BW:30:GLU:N	2.25	0.52
55:BY:13:VAL:HG12	55:BY:14:LEU:N	2.24	0.52
56:BZ:3:TYR:O	56:BZ:57:ILE:HG23	2.09	0.52
56:BZ:27:VAL:O	56:BZ:87:ASP:HA	2.10	0.52
1:CA:1495:U:H2'	1:CA:1496:C:H6	1.74	0.52
1:CA:167:G:O2'	1:CA:168:G:H5'	2.10	0.52
1:CA:269:C:H2'	1:CA:270:A:H8	1.75	0.52
1:CA:491:G:H2'	1:CA:492:G:H8	1.74	0.52
2:CB:175:ARG:O	2:CB:176:GLU:C	2.47	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:212:GLN:HE22	2:CB:216:SER:CB	2.22	0.52
3:CC:178:LEU:HD22	3:CC:178:LEU:N	2.25	0.52
3:CC:47:LEU:CD2	3:CC:52:LEU:HD13	2.40	0.52
7:CG:143:ARG:NH1	7:CG:143:ARG:HB2	2.23	0.52
1:CA:1343:G:H4'	9:CI:122:ALA:O	2.10	0.52
12:CL:8:ASN:O	12:CL:11:VAL:HB	2.09	0.52
12:CL:119:LYS:C	12:CL:120:TYR:CD1	2.83	0.52
12:CL:90:VAL:C	12:CL:92:ASP:H	2.13	0.52
13:CM:108:ARG:NH2	13:CM:114:ARG:HG2	2.24	0.52
15:CO:17:ARG:NH1	15:CO:77:ARG:NH1	2.57	0.52
1:CA:237:C:C4'	17:CQ:25:ARG:HH12	2.18	0.52
23:CW:38:A:H5'	23:CW:39:A:OP2	2.10	0.52
26:D0:72:ARG:CD	26:D0:75:LEU:HD13	2.38	0.52
35:DA:1819:A:H5''	38:DD:161:THR:CG2	2.40	0.52
35:DA:2639:A:H3'	35:DA:2640:G:H5'	1.92	0.52
35:DA:2774:C:H2'	35:DA:2775:A:H8	1.75	0.52
35:DA:2821:A:OP2	48:DR:2:ARG:NH2	2.42	0.52
35:DA:309:G:H4'	55:DY:18:GLY:HA3	1.92	0.52
35:DA:338:G:H2'	35:DA:339:U:C6	2.44	0.52
35:DA:60:G:N2	35:DA:74:A:H2'	2.25	0.52
35:DA:711:G:O2'	35:DA:712:G:H5'	2.10	0.52
36:DB:60:C:C2	36:DB:61:G:C8	2.98	0.52
35:DA:782:A:C2	38:DD:226:MET:CE	2.93	0.52
39:DE:119:ARG:NH1	39:DE:159:HIS:O	2.42	0.52
39:DE:128:SER:O	39:DE:129:HIS:CB	2.56	0.52
39:DE:60:ASN:O	39:DE:63:LEU:N	2.43	0.52
35:DA:2632:A:N3	39:DE:61:ARG:NH1	2.58	0.52
40:DF:108:LYS:O	40:DF:112:MET:SD	2.68	0.52
40:DF:185:ASP:O	40:DF:189:THR:HG23	2.09	0.52
42:DH:130:ARG:HH11	42:DH:130:ARG:HB2	1.75	0.52
45:DO:34:THR:O	45:DO:37:ASP:OD2	2.27	0.52
48:DR:2:ARG:HE	48:DR:5:LYS:HZ2	1.57	0.52
50:DT:50:ILE:H	50:DT:50:ILE:HD12	1.74	0.52
45:DO:77:ILE:HD11	50:DT:72:VAL:HG13	1.92	0.52
51:DU:18:LEU:HD21	51:DU:22:LYS:HE2	1.91	0.52
52:DV:5:VAL:HG23	52:DV:37:VAL:H	1.75	0.52
55:DY:96:ILE:HG13	55:DY:99:CYS:HB2	1.90	0.52
1:AA:1208:C:H2'	1:AA:1209:C:H6	1.75	0.52
1:AA:1354:C:O2'	1:AA:1355:G:H5'	2.10	0.52
1:AA:1504:G:OP1	1:AA:1507:A:H4'	2.09	0.52
1:AA:44:G:C2	1:AA:45:U:H1'	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:764:C:H2'	1:AA:765:G:O4'	2.10	0.52
1:AA:806:C:O2'	1:AA:807:A:H5'	2.11	0.52
1:AA:894:G:H2'	1:AA:895:G:H8	1.75	0.52
2:AB:74:LYS:HG3	2:AB:77:ALA:HB3	1.92	0.52
3:AC:141:VAL:HG11	3:AC:202:ILE:HD12	1.92	0.52
4:AD:13:ARG:HD3	4:AD:39:PRO:C	2.30	0.52
4:AD:61:LYS:HG2	4:AD:75:PHE:HE2	1.75	0.52
5:AE:79:GLU:N	5:AE:79:GLU:OE1	2.40	0.52
7:AG:25:ALA:HA	7:AG:28:ASN:HD22	1.75	0.52
7:AG:49:ILE:HG22	7:AG:53:LYS:HG3	1.92	0.52
9:AI:65:VAL:CG2	9:AI:66:ARG:N	2.72	0.52
1:AA:692:U:H5	11:AK:26:ASN:ND2	2.08	0.52
12:AL:20:LYS:H	12:AL:20:LYS:HD3	1.75	0.52
12:AL:21:LYS:CD	12:AL:21:LYS:H	2.10	0.52
12:AL:85:ILE:HD11	12:AL:98:TYR:HB2	1.90	0.52
15:AO:39:LEU:HD12	15:AO:56:LEU:CD1	2.40	0.52
19:AS:53:ASN:N	19:AS:53:ASN:HD22	2.06	0.52
19:AS:36:ARG:CZ	19:AS:75:ALA:HB3	2.39	0.52
20:AT:14:LYS:HE3	20:AT:18:GLN:NE2	2.25	0.52
20:AT:41:ILE:O	20:AT:43:LEU:N	2.43	0.52
27:B1:8:SER:HA	35:BA:1365:A:OP2	2.10	0.52
33:B7:11:LYS:O	33:B7:14:LYS:N	2.41	0.52
33:B7:8:ASN:HD22	33:B7:8:ASN:C	2.12	0.52
35:BA:1014:U:H2'	35:BA:1015:G:O4'	2.09	0.52
35:BA:104:U:H6	35:BA:104:U:O5'	1.92	0.52
35:BA:1195:G:H2'	35:BA:1196:C:H6	1.74	0.52
35:BA:1227:G:OP2	51:BU:16:LYS:HE3	2.10	0.52
35:BA:1289:C:H2'	35:BA:1290:C:H6	1.73	0.52
35:BA:1779:U:C2	35:BA:1783:A:N7	2.78	0.52
35:BA:1990:C:H2'	35:BA:1991:U:O4'	2.10	0.52
35:BA:2785:C:H2'	35:BA:2786:U:C6	2.45	0.52
35:BA:335:C:H4'	55:BY:73:ARG:CZ	2.40	0.52
35:BA:836:G:H2'	35:BA:837:C:H6	1.72	0.52
35:BA:910:A:H2'	35:BA:911:A:C8	2.44	0.52
35:BA:970:C:H2'	35:BA:971:C:C6	2.44	0.52
36:BB:45:A:N3	36:BB:45:A:H2'	2.23	0.52
38:BD:112:GLN:N	38:BD:112:GLN:OE1	2.43	0.52
39:BE:2:LYS:CE	39:BE:95:ILE:HG22	2.39	0.52
39:BE:59:VAL:CG1	39:BE:63:LEU:HG	2.40	0.52
35:BA:2810:A:H2'	39:BE:61:ARG:NH2	2.25	0.52
42:BH:147:ASN:O	42:BH:150:ALA:HB3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BH:35:VAL:HG12	42:BH:35:VAL:O	2.10	0.52
42:BH:40:GLU:O	42:BH:41:MET:HB2	2.10	0.52
45:BO:87:ILE:CG2	45:BO:88:ASN:N	2.72	0.52
46:BP:74:GLU:HA	46:BP:74:GLU:OE2	2.10	0.52
39:BE:110:GLY:O	48:BR:2:ARG:NH2	2.43	0.52
48:BR:28:LEU:C	48:BR:30:THR:H	2.13	0.52
49:BS:13:ARG:H	49:BS:13:ARG:CD	1.99	0.52
51:BU:112:ARG:HH11	51:BU:112:ARG:HG2	1.75	0.52
54:BX:29:TRP:HE3	54:BX:29:TRP:HA	1.75	0.52
54:BX:85:PRO:O	54:BX:86:GLY:C	2.48	0.52
56:BZ:146:ILE:O	56:BZ:148:ASP:N	2.43	0.52
1:CA:1300:G:O2'	1:CA:1301:U:O5'	2.26	0.52
1:CA:328:C:C2'	1:CA:328:C:O2	2.58	0.52
1:CA:358:U:O2'	1:CA:359:U:H5'	2.10	0.52
1:CA:444:C:H2'	1:CA:445:G:C8	2.41	0.52
1:CA:806:C:O2'	1:CA:807:A:H5'	2.09	0.52
2:CB:163:PHE:O	2:CB:164:VAL:HG23	2.10	0.52
2:CB:82:ARG:HB2	2:CB:94:ASN:ND2	2.24	0.52
3:CC:141:VAL:HG11	3:CC:202:ILE:HD12	1.91	0.52
4:CD:190:ASP:O	4:CD:191:ARG:C	2.48	0.52
10:CJ:7:LYS:O	10:CJ:96:ILE:HA	2.10	0.52
11:CK:60:ALA:O	11:CK:61:ALA:C	2.49	0.52
15:CO:81:LEU:O	15:CO:84:LYS:HB2	2.09	0.52
16:CP:72:ARG:C	16:CP:74:LEU:N	2.63	0.52
1:CA:237:C:H5''	17:CQ:25:ARG:NH1	2.25	0.52
22:CV:35:A:H2'	22:CV:36:A:O4'	2.10	0.52
25:CY:117:ALA:O	25:CY:120:GLN:HB3	2.09	0.52
27:D1:23:LYS:C	27:D1:37:ILE:HG12	2.30	0.52
27:D1:59:THR:HG23	27:D1:59:THR:O	2.09	0.52
28:D2:22:GLU:C	28:D2:24:LEU:H	2.13	0.52
29:D3:41:PRO:HD3	29:D3:44:ARG:CZ	2.40	0.52
29:D3:54:VAL:HG12	29:D3:55:ARG:N	2.24	0.52
34:D8:60:LEU:N	34:D8:60:LEU:CD2	2.72	0.52
35:DA:109:G:O2'	35:DA:110:G:H5'	2.10	0.52
35:DA:1335:U:H2'	35:DA:1336:A:H8	1.74	0.52
35:DA:188:G:H1'	35:DA:1365:A:N1	2.25	0.52
35:DA:1667:G:OP1	45:DO:7:TYR:HB2	2.10	0.52
35:DA:1695:G:H2'	35:DA:1696:G:C5'	2.40	0.52
35:DA:174:C:O2	35:DA:174:C:H2'	2.09	0.52
35:DA:1803:A:C8	35:DA:1804:C:C5	2.98	0.52
35:DA:197:A:H5'	35:DA:197:A:C8	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:251:A:H2'	35:DA:252:G:O4'	2.09	0.52
35:DA:2768:C:O2'	35:DA:2769:C:H5'	2.10	0.52
35:DA:2875:C:C4'	50:DT:5:ALA:HB2	2.39	0.52
35:DA:587:C:C5	35:DA:671:C:H1'	2.44	0.52
38:DD:65:ILE:HD11	38:DD:67:PHE:CD1	2.44	0.52
40:DF:154:VAL:HB	40:DF:173:VAL:HG22	1.90	0.52
40:DF:7:TYR:HB3	40:DF:16:GLY:O	2.09	0.52
41:DG:114:ILE:HD12	41:DG:117:PHE:CD1	2.45	0.52
41:DG:16:ARG:HB3	41:DG:17:PRO:CD	2.40	0.52
41:DG:46:ALA:HA	41:DG:51:ARG:CD	2.40	0.52
41:DG:86:MET:N	41:DG:87:PRO:CD	2.73	0.52
45:DO:37:ASP:O	45:DO:61:VAL:HA	2.10	0.52
46:DP:100:LEU:H	46:DP:100:LEU:CD2	2.23	0.52
46:DP:56:SER:C	46:DP:58:THR:H	2.13	0.52
47:DQ:20:ALA:HB2	47:DQ:99:PRO:CD	2.40	0.52
47:DQ:59:ARG:HH11	47:DQ:59:ARG:HG3	1.75	0.52
51:DU:59:ARG:O	51:DU:60:LEU:C	2.48	0.52
54:DX:25:LYS:O	54:DX:26:TYR:O	2.28	0.52
55:DY:101:LYS:CG	55:DY:102:CYS:N	2.72	0.52
55:DY:14:LEU:CD1	55:DY:15:VAL:H	2.19	0.52
1:AA:1333:A:C2	1:AA:1334:G:H1'	2.45	0.51
1:AA:385:C:H2'	1:AA:386:C:H6	1.75	0.51
1:AA:498:U:C2'	1:AA:499:A:H5'	2.40	0.51
1:AA:594:G:O2'	1:AA:595:G:H5'	2.09	0.51
3:AC:109:PRO:O	3:AC:115:LEU:HD12	2.10	0.51
1:AA:1112:C:C4	3:AC:178:LEU:HD23	2.45	0.51
5:AE:127:ASN:O	5:AE:131:ILE:HG12	2.10	0.51
5:AE:31:LEU:CD1	5:AE:129:ILE:HA	2.40	0.51
6:AF:39:LYS:HG2	6:AF:40:VAL:N	2.23	0.51
10:AJ:78:ASN:C	10:AJ:80:LYS:N	2.63	0.51
11:AK:102:GLY:O	11:AK:103:LEU:C	2.47	0.51
11:AK:58:PRO:HD3	11:AK:89:ALA:CB	2.40	0.51
11:AK:21:ILE:CB	11:AK:84:VAL:HG12	2.39	0.51
13:AM:3:ARG:HA	13:AM:9:ILE:CG1	2.37	0.51
13:AM:86:CYS:O	13:AM:89:GLY:N	2.44	0.51
15:AO:75:PRO:O	15:AO:79:ARG:HG3	2.10	0.51
15:AO:81:LEU:O	15:AO:84:LYS:HB2	2.11	0.51
16:AP:60:LEU:C	16:AP:62:VAL:N	2.64	0.51
17:AQ:23:VAL:O	17:AQ:39:SER:HB2	2.09	0.51
17:AQ:43:LEU:HD12	17:AQ:68:ARG:HG3	1.91	0.51
18:AR:25:THR:O	18:AR:26:LEU:HG	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:AY:105:PRO:O	25:AY:106:LEU:HD23	2.10	0.51
31:B5:30:LEU:HD23	31:B5:41:PRO:CA	2.40	0.51
33:B7:16:HIS:CE1	35:BA:684:G:OP1	2.63	0.51
35:BA:1300:U:O2'	35:BA:1301:A:P	2.68	0.51
35:BA:1710:C:O2'	35:BA:1711:C:H5'	2.10	0.51
35:BA:2280:G:N3	35:BA:2388:A:H2	2.08	0.51
35:BA:2553:G:H2'	35:BA:2554:U:H4'	1.92	0.51
35:BA:2870:C:O2'	35:BA:2871:C:H5'	2.10	0.51
35:BA:2875:C:C4'	50:BT:5:ALA:HB2	2.39	0.51
36:BB:52:A:HO2'	36:BB:53:A:H8	1.52	0.51
41:BG:16:ARG:HH11	41:BG:16:ARG:HG3	1.76	0.51
42:BH:163:TYR:CD1	42:BH:163:TYR:N	2.78	0.51
44:BN:17:ASP:OD2	44:BN:55:VAL:O	2.27	0.51
35:BA:1190:G:C4'	46:BP:35:HIS:HB3	2.40	0.51
49:BS:11:LYS:N	49:BS:11:LYS:HD3	2.24	0.51
53:BW:59:VAL:O	53:BW:61:ASN:N	2.43	0.51
54:BX:55:ASN:C	54:BX:77:LYS:CG	2.75	0.51
54:BX:83:VAL:O	54:BX:84:ALA:CB	2.58	0.51
55:BY:10:GLY:HA2	55:BY:27:VAL:CG1	2.31	0.51
55:BY:39:VAL:O	55:BY:40:GLU:HG2	2.10	0.51
1:CA:1261:A:H62	1:CA:1274:G:N2	2.07	0.51
1:CA:1332:A:H2'	1:CA:1333:A:H8	1.75	0.51
1:CA:376:G:H2'	1:CA:377:G:C8	2.46	0.51
1:CA:972:C:C5'	10:CJ:57:LYS:HG3	2.40	0.51
3:CC:3:ASN:O	3:CC:4:LYS:C	2.48	0.51
4:CD:110:PHE:HD1	4:CD:110:PHE:H	1.58	0.51
4:CD:110:PHE:HD1	4:CD:110:PHE:N	2.08	0.51
5:CE:127:ASN:O	5:CE:131:ILE:HG12	2.10	0.51
10:CJ:34:VAL:HG12	10:CJ:35:SER:H	1.75	0.51
11:CK:103:LEU:N	11:CK:103:LEU:CD2	2.69	0.51
16:CP:22:THR:CG2	16:CP:32:TYR:HA	2.39	0.51
25:CY:16:LYS:C	25:CY:18:LEU:N	2.63	0.51
27:D1:72:GLU:O	27:D1:76:ARG:NH2	2.43	0.51
33:D7:29:LYS:NZ	33:D7:32:LYS:HZ2	2.08	0.51
34:D8:49:VAL:CB	34:D8:53:PRO:HD3	2.40	0.51
34:D8:51:ALA:HA	34:D8:54:GLU:OE1	2.10	0.51
35:DA:1301:A:O2'	35:DA:1302:A:C2'	2.56	0.51
35:DA:1531:C:H3'	35:DA:1532:C:O4'	2.10	0.51
35:DA:2359:C:H6	35:DA:2359:C:O5'	1.93	0.51
35:DA:2393:A:O2'	35:DA:2394:C:H5'	2.11	0.51
35:DA:2425:A:O4'	35:DA:2427:C:C6	2.63	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2467:C:O2'	35:DA:2468:G:H5'	2.11	0.51
35:DA:2511:U:O3'	39:DE:123:ALA:HB3	2.10	0.51
35:DA:2846:G:H2'	35:DA:2847:U:C6	2.44	0.51
35:DA:290:G:C2'	35:DA:291:C:H5'	2.40	0.51
35:DA:554:U:C2'	35:DA:555:U:H5'	2.41	0.51
38:DD:209:ALA:O	38:DD:210:GLY:O	2.28	0.51
35:DA:1658:C:OP1	39:DE:132:HIS:O	2.28	0.51
42:DH:117:PRO:HB3	42:DH:123:PHE:HD1	1.75	0.51
42:DH:148:ILE:N	42:DH:148:ILE:HD13	2.25	0.51
46:DP:97:PRO:C	46:DP:99:LEU:H	2.13	0.51
48:DR:104:ARG:O	48:DR:106:GLY:N	2.43	0.51
48:DR:16:HIS:O	48:DR:17:ARG:C	2.49	0.51
49:DS:83:LYS:O	49:DS:85:VAL:HG22	2.10	0.51
50:DT:108:ARG:O	50:DT:112:ARG:HG3	2.10	0.51
50:DT:91:ARG:HB3	50:DT:115:ARG:O	2.10	0.51
50:DT:28:VAL:HB	50:DT:88:ILE:HG12	1.91	0.51
50:DT:32:TYR:O	50:DT:33:LYS:CB	2.57	0.51
50:DT:40:THR:O	50:DT:41:ARG:CB	2.48	0.51
51:DU:38:THR:O	51:DU:41:ALA:HB3	2.09	0.51
52:DV:69:LYS:HB2	52:DV:93:GLU:OE2	2.10	0.51
53:DW:17:VAL:C	53:DW:19:LEU:N	2.63	0.51
55:DY:68:HIS:HB3	55:DY:71:LYS:HZ3	1.75	0.51
56:DZ:18:LEU:HG	56:DZ:23:LYS:CB	2.40	0.51
1:AA:1148:U:OP1	9:AI:7:THR:HG21	2.10	0.51
1:AA:1392:G:O2'	1:AA:1393:U:H5'	2.10	0.51
1:AA:199:G:H2'	1:AA:200:G:H8	1.75	0.51
1:AA:23:C:O2'	1:AA:24:U:H5'	2.10	0.51
1:AA:895:G:H2'	1:AA:896:C:C6	2.45	0.51
2:AB:175:ARG:O	2:AB:177:ALA:N	2.43	0.51
2:AB:41:ILE:HD12	2:AB:41:ILE:N	2.24	0.51
3:AC:109:PRO:HG2	3:AC:110:ASN:H	1.73	0.51
3:AC:24:ALA:HB1	3:AC:28:GLN:O	2.10	0.51
4:AD:202:LEU:O	4:AD:205:GLU:N	2.42	0.51
4:AD:87:GLY:C	4:AD:89:THR:H	2.12	0.51
6:AF:100:ASN:O	6:AF:101:ALA:O	2.27	0.51
6:AF:12:PRO:HG3	6:AF:57:GLN:O	2.10	0.51
11:AK:22:HIS:C	11:AK:28:THR:HG23	2.30	0.51
12:AL:60:LEU:O	12:AL:62:SER:N	2.44	0.51
13:AM:35:GLU:HG3	13:AM:36:LYS:N	2.24	0.51
1:AA:267:C:P	17:AQ:67:LYS:HB2	2.50	0.51
17:AQ:56:VAL:CG2	17:AQ:78:GLU:HG3	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AT:32:ALA:C	20:AT:36:LEU:HD23	2.31	0.51
27:B1:10:LYS:C	27:B1:13:ILE:HG22	2.28	0.51
27:B1:47:GLN:HE21	27:B1:64:ALA:HB2	1.74	0.51
27:B1:87:PRO:HG2	27:B1:88:LYS:N	2.26	0.51
35:BA:1112:G:C1'	35:BA:1113:U:OP1	2.57	0.51
35:BA:1248:G:H2'	51:BU:2:PRO:O	2.11	0.51
35:BA:1434:A:O2'	35:BA:1435:G:H5'	2.10	0.51
35:BA:1531:C:H3'	35:BA:1532:C:O4'	2.10	0.51
35:BA:1931:U:O2'	35:BA:1932:A:H5'	2.11	0.51
35:BA:2007:C:H6	35:BA:2007:C:O5'	1.93	0.51
35:BA:2248:C:O2'	35:BA:2249:U:H5'	2.11	0.51
35:BA:52:A:C2'	35:BA:53:A:H5'	2.40	0.51
35:BA:824:A:H2'	35:BA:825:C:C6	2.45	0.51
35:BA:959:A:H2'	35:BA:960:A:C8	2.44	0.51
35:BA:998:C:N4	35:BA:1158:C:H42	2.07	0.51
38:BD:131:LEU:N	38:BD:131:LEU:HD12	2.26	0.51
38:BD:92:ILE:HA	38:BD:107:ALA:HB2	1.92	0.51
41:BG:24:GLY:O	41:BG:25:TYR:C	2.48	0.51
42:BH:117:PRO:HB3	42:BH:123:PHE:HD1	1.75	0.51
42:BH:85:LYS:HZ3	42:BH:145:ALA:HA	1.74	0.51
42:BH:89:ILE:CG1	42:BH:90:LYS:H	2.22	0.51
44:BN:52:VAL:O	44:BN:120:LEU:HD22	2.10	0.51
44:BN:78:TYR:CG	44:BN:79:PRO:HD3	2.45	0.51
45:BO:47:ILE:HG23	45:BO:48:PRO:CD	2.40	0.51
35:BA:244:A:O2'	46:BP:73:GLY:HA3	2.10	0.51
47:BQ:34:LEU:HD22	47:BQ:121:ALA:HB3	1.92	0.51
50:BT:43:GLN:HG2	50:BT:44:ASP:O	2.11	0.51
51:BU:92:ARG:NH1	52:BV:11:GLN:O	2.42	0.51
54:BX:29:TRP:CE3	54:BX:29:TRP:HA	2.44	0.51
55:BY:28:LYS:CD	55:BY:37:VAL:HG12	2.40	0.51
56:BZ:156:LYS:O	56:BZ:156:LYS:HG2	2.10	0.51
1:CA:1170:A:H2'	1:CA:1171:G:H5'	1.91	0.51
1:CA:1388:C:H2'	1:CA:1389:C:H6	1.70	0.51
1:CA:1530:G:OP1	1:CA:1530:G:H4'	2.10	0.51
1:CA:325:A:N6	1:CA:326:G:C6	2.78	0.51
1:CA:702:A:H3'	1:CA:703:G:C5'	2.39	0.51
1:CA:788:U:O2'	1:CA:789:U:H5'	2.10	0.51
2:CB:28:PHE:HD1	2:CB:28:PHE:O	1.94	0.51
3:CC:141:VAL:HG11	3:CC:202:ILE:CD1	2.40	0.51
3:CC:40:ARG:HB3	3:CC:44:GLU:OE2	2.10	0.51
4:CD:10:ARG:C	4:CD:13:ARG:HB3	2.30	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:114:THR:HG23	8:CH:116:LYS:O	2.10	0.51
17:CQ:29:HIS:CA	17:CQ:36:ILE:HD11	2.40	0.51
1:CA:235:C:H5'	17:CQ:70:ARG:HG2	1.92	0.51
19:CS:50:ALA:HB1	19:CS:57:HIS:HB3	1.91	0.51
28:D2:30:ARG:O	28:D2:34:GLU:HB3	2.10	0.51
28:D2:48:HIS:CE1	35:DA:75:G:H4'	2.45	0.51
29:D3:22:ALA:CB	29:D3:46:ASN:HD22	2.23	0.51
35:DA:1153:C:N4	35:DA:1154:G:C2	2.79	0.51
35:DA:149:A:H2'	35:DA:150:C:O4'	2.09	0.51
35:DA:1600:C:C2'	35:DA:1601:G:H5'	2.40	0.51
35:DA:1771:C:H1'	35:DA:1786:A:H8	1.74	0.51
35:DA:2056:G:N2	35:DA:2057:A:C1'	2.72	0.51
26:D0:41:ARG:HH21	35:DA:2387:U:H1'	1.73	0.51
35:DA:2413:G:H21	46:DP:70:GLN:NE2	2.08	0.51
35:DA:2562:U:H2'	35:DA:2563:U:C5'	2.40	0.51
35:DA:271(F):C:H2'	35:DA:271(G):C:H6	1.74	0.51
35:DA:271(U):G:O2'	35:DA:271(V):G:H5'	2.10	0.51
35:DA:2759:G:H5'	35:DA:2759:G:C8	2.45	0.51
35:DA:289:A:H3'	35:DA:290:G:H8	1.75	0.51
35:DA:589:C:O2'	35:DA:590:A:H5'	2.10	0.51
36:DB:87:G:H2'	36:DB:88:C:H5''	1.91	0.51
42:DH:147:ASN:O	42:DH:150:ALA:HB3	2.10	0.51
43:DI:4:ILE:O	43:DI:36:ALA:HB1	2.10	0.51
43:DI:5:LEU:O	43:DI:6:LEU:HG	2.11	0.51
44:DN:97:ARG:HB3	44:DN:101:HIS:HD2	1.76	0.51
35:DA:1246:A:OP1	46:DP:18:ARG:HG3	2.11	0.51
47:DQ:66:ILE:HG22	47:DQ:104:PHE:HE2	1.73	0.51
47:DQ:20:ALA:O	47:DQ:23:GLY:N	2.34	0.51
48:DR:10:LEU:HB3	48:DR:17:ARG:NE	2.24	0.51
50:DT:100:TYR:C	50:DT:102:ILE:N	2.63	0.51
50:DT:28:VAL:CG2	50:DT:46:GLU:HA	2.40	0.51
39:DE:9:VAL:HG23	50:DT:8:LYS:HB2	1.92	0.51
51:DU:92:ARG:NH2	52:DV:10:LYS:HA	2.25	0.51
54:DX:7:VAL:HB	54:DX:8:ILE:HD12	1.92	0.51
56:DZ:45:ASP:O	56:DZ:46:LYS:C	2.49	0.51
1:AA:1051:C:O2'	1:AA:1052:U:H5'	2.11	0.51
1:AA:1332:A:H2'	1:AA:1333:A:H8	1.75	0.51
1:AA:994:A:H2	14:AN:4:LYS:HG3	1.74	0.51
3:AC:141:VAL:HG11	3:AC:202:ILE:CD1	2.41	0.51
3:AC:172:ARG:HH11	3:AC:172:ARG:HB3	1.74	0.51
4:AD:196:LEU:HD12	4:AD:196:LEU:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AE:102:ALA:HB1	5:AE:106:PRO:HG2	1.92	0.51
5:AE:71:LEU:HD11	5:AE:114:GLY:C	2.30	0.51
7:AG:145:ALA:O	7:AG:146:GLU:HB2	2.09	0.51
8:AH:64:LYS:O	8:AH:79:VAL:CG2	2.59	0.51
8:AH:83:ILE:HD13	8:AH:137:VAL:CG2	2.34	0.51
10:AJ:22:LYS:NZ	10:AJ:23:ILE:HG12	2.25	0.51
12:AL:38:THR:HG23	12:AL:39:VAL:N	2.26	0.51
13:AM:68:GLY:O	13:AM:71:ARG:N	2.43	0.51
14:AN:40:CYS:SG	14:AN:43:CYS:N	2.75	0.51
14:AN:37:PHE:HE1	14:AN:53:LEU:HD22	1.75	0.51
15:AO:66:LEU:H	15:AO:66:LEU:HD12	1.75	0.51
1:AA:835:U:OP1	18:AR:61:LYS:HB2	2.10	0.51
23:AW:29:C:H2'	23:AW:30:G:C8	2.43	0.51
29:B3:22:ALA:CB	29:B3:46:ASN:HD22	2.23	0.51
30:B4:29:PRO:C	30:B4:31:ILE:N	2.63	0.51
32:B6:10:LEU:HD22	32:B6:10:LEU:H	1.76	0.51
35:BA:1213:A:N3	35:BA:1238:G:H1'	2.25	0.51
35:BA:1515:G:O2'	35:BA:1516:C:H5'	2.11	0.51
35:BA:1572:A:O2'	35:BA:1573:G:H5'	2.10	0.51
35:BA:1612:C:H2'	35:BA:1613:G:O5'	2.11	0.51
35:BA:2079:U:H2'	35:BA:2080:G:C8	2.45	0.51
35:BA:27:G:H1'	35:BA:513:A:H62	1.76	0.51
35:BA:2821:A:OP2	35:BA:2822:G:OP2	2.28	0.51
35:BA:402:A:O2'	35:BA:403:U:H5'	2.10	0.51
35:BA:675:A:O2'	35:BA:676:A:H5'	2.09	0.51
38:BD:111:LEU:HD13	38:BD:112:GLN:N	2.25	0.51
38:BD:132:PRO:O	38:BD:136:ILE:HG13	2.10	0.51
45:BO:105:GLU:HA	45:BO:108:GLU:HG3	1.92	0.51
46:BP:21:ARG:NH1	46:BP:21:ARG:HG3	2.24	0.51
46:BP:70:GLN:HA	46:BP:70:GLN:OE1	2.10	0.51
47:BQ:53:ALA:O	47:BQ:56:ARG:HB3	2.10	0.51
48:BR:2:ARG:NE	48:BR:5:LYS:HE3	2.25	0.51
53:BW:48:ALA:O	53:BW:51:LEU:N	2.43	0.51
55:BY:74:PRO:HG2	55:BY:80:GLY:O	2.10	0.51
1:CA:119:A:H4'	1:CA:120:A:O5'	2.10	0.51
1:CA:1397:C:H5''	1:CA:1398:A:O5'	2.10	0.51
1:CA:390:C:O3'	16:CP:28:ARG:NH2	2.43	0.51
1:CA:651:C:H2'	1:CA:652:U:C6	2.45	0.51
1:CA:794:A:H2'	1:CA:795:C:C6	2.45	0.51
1:CA:865:A:H2'	1:CA:866:C:H6	1.75	0.51
4:CD:120:LEU:HA	4:CD:125:HIS:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:922:G:H4'	5:CE:20:GLN:HG2	1.91	0.51
7:CG:106:GLN:HA	7:CG:109:ASN:HD22	1.76	0.51
7:CG:104:LEU:O	7:CG:107:ALA:HB3	2.10	0.51
9:CI:24:GLY:HA2	9:CI:59:PHE:O	2.11	0.51
9:CI:6:GLY:HA3	9:CI:84:ALA:N	2.26	0.51
13:CM:23:TYR:HE1	13:CM:71:ARG:CB	2.19	0.51
14:CN:3:ARG:NH1	14:CN:3:ARG:HB3	2.26	0.51
18:CR:32:ARG:HA	18:CR:69:THR:HG21	1.92	0.51
25:CY:108:GLU:HG2	25:CY:111:ARG:NH2	2.25	0.51
25:CY:129:ILE:O	25:CY:132:ILE:HG13	2.10	0.51
25:CY:2:THR:O	25:CY:3:LEU:C	2.49	0.51
25:CY:38:LEU:C	25:CY:40:HIS:H	2.14	0.51
27:D1:47:GLN:OE1	27:D1:47:GLN:HA	2.10	0.51
28:D2:23:LYS:HB2	54:DX:5:TYR:CE1	2.46	0.51
34:D8:55:ALA:O	34:D8:59:LYS:NZ	2.36	0.51
34:D8:56:GLU:O	34:D8:58:ILE:N	2.43	0.51
35:DA:1277:G:O2'	35:DA:1278:A:H5'	2.11	0.51
35:DA:2550:G:H2'	35:DA:2551:C:H6	1.75	0.51
35:DA:2692:C:H2'	35:DA:2693:A:H8	1.75	0.51
35:DA:2886:G:N2	35:DA:2887:U:C2	2.79	0.51
35:DA:514:A:H2'	35:DA:515:A:H8	1.74	0.51
38:DD:14:ARG:HH11	38:DD:14:ARG:CG	2.23	0.51
38:DD:158:ALA:O	38:DD:159:ALA:C	2.49	0.51
39:DE:49:LEU:HD23	39:DE:81:ILE:HG12	1.91	0.51
39:DE:78:LEU:O	39:DE:79:ARG:HD2	2.10	0.51
40:DF:39:TRP:HA	40:DF:99:TYR:CE1	2.45	0.51
40:DF:65:TRP:O	40:DF:67:GLN:N	2.44	0.51
41:DG:108:ASN:C	41:DG:112:PRO:HD2	2.31	0.51
41:DG:130:ASN:HA	41:DG:161:THR:OG1	2.10	0.51
41:DG:21:ARG:HD3	41:DG:21:ARG:C	2.31	0.51
42:DH:85:LYS:NZ	42:DH:145:ALA:HA	2.25	0.51
43:DI:58:LEU:HD23	43:DI:58:LEU:C	2.30	0.51
45:DO:104:ARG:C	45:DO:106:LEU:N	2.61	0.51
45:DO:35:VAL:CG2	45:DO:69:ILE:HG12	2.39	0.51
49:DS:62:LYS:O	49:DS:65:VAL:N	2.43	0.51
53:DW:66:GLU:HG2	53:DW:66:GLU:O	2.10	0.51
54:DX:65:ARG:CZ	54:DX:66:LEU:N	2.68	0.51
55:DY:54:LYS:O	55:DY:55:TYR:O	2.29	0.51
56:DZ:59:LEU:H	56:DZ:59:LEU:CD2	2.23	0.51
1:AA:1298:C:H1'	1:AA:1299:A:C2	2.45	0.51
1:AA:1349:A:H2'	1:AA:1350:A:H8	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1429:C:O2'	1:AA:1430:C:H5'	2.10	0.51
1:AA:969:A:H2'	1:AA:970:C:O4'	2.10	0.51
2:AB:19:HIS:CA	2:AB:39:ILE:HD13	2.39	0.51
4:AD:22:LYS:HB3	4:AD:22:LYS:NZ	2.25	0.51
6:AF:42:GLU:C	6:AF:44:GLY:N	2.64	0.51
9:AI:13:ALA:HA	9:AI:67:GLY:C	2.30	0.51
9:AI:3:GLN:CG	9:AI:20:ARG:HH21	2.23	0.51
9:AI:65:VAL:HG22	9:AI:66:ARG:H	1.74	0.51
14:AN:29:ARG:HG3	14:AN:29:ARG:HH11	1.75	0.51
17:AQ:29:HIS:CE1	17:AQ:31:LEU:HB3	2.45	0.51
24:AX:20:A:O2'	24:AX:21:G:H5'	2.10	0.51
27:B1:47:GLN:NE2	27:B1:48:LYS:O	2.43	0.51
29:B3:40:THR:HG23	29:B3:43:ILE:H	1.75	0.51
32:B6:15:GLU:O	32:B6:15:GLU:CG	2.58	0.51
35:BA:1225:G:OP1	52:BV:88:ARG:HD2	2.10	0.51
35:BA:2119:A:C3'	35:BA:2120:G:C5'	2.89	0.51
35:BA:2379:G:H2'	35:BA:2380:C:C6	2.46	0.51
35:BA:2407:G:C2	35:BA:2408:U:C4	2.99	0.51
35:BA:2420:C:O2'	35:BA:2421:G:H5'	2.10	0.51
35:BA:2718:G:H2'	35:BA:2719:G:H8	1.76	0.51
35:BA:391:G:O2'	35:BA:392:C:H5'	2.11	0.51
35:BA:836:G:C5	35:BA:837:C:C4	2.98	0.51
35:BA:873:G:H1	35:BA:904:C:H42	1.58	0.51
37:BC:214:VAL:O	37:BC:216:THR:N	2.43	0.51
38:BD:16:MET:HG3	38:BD:206:LEU:O	2.10	0.51
35:BA:782:A:C2	38:BD:226:MET:CE	2.92	0.51
39:BE:47:VAL:CG2	39:BE:84:PHE:O	2.59	0.51
40:BF:21:ALA:O	40:BF:23:ASP:N	2.44	0.51
41:BG:77:ILE:HD12	41:BG:82:LEU:O	2.10	0.51
49:BS:29:PHE:HB3	49:BS:36:TYR:HB2	1.92	0.51
49:BS:32:LEU:O	49:BS:33:LYS:HB2	2.09	0.51
49:BS:62:LYS:O	49:BS:65:VAL:N	2.44	0.51
52:BV:63:GLY:O	52:BV:64:HIS:HB3	2.11	0.51
53:BW:20:VAL:O	53:BW:23:LEU:HB2	2.10	0.51
56:BZ:121:HIS:C	56:BZ:123:ASP:N	2.64	0.51
1:CA:1255:G:H2'	1:CA:1255:G:N3	2.24	0.51
1:CA:139:G:O2'	1:CA:140:A:H5'	2.11	0.51
1:CA:1405:G:H21	1:CA:1517:G:H22	1.55	0.51
1:CA:522:C:N4	1:CA:528:C:N4	2.56	0.51
1:CA:997:U:H2'	1:CA:998:G:C8	2.45	0.51
4:CD:105:VAL:CG2	4:CD:126:ILE:HD13	2.37	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:13:ARG:HD3	4:CD:39:PRO:C	2.30	0.51
4:CD:13:ARG:HG2	4:CD:14:ARG:H	1.74	0.51
5:CE:12:LEU:O	5:CE:12:LEU:HD22	2.10	0.51
5:CE:89:ILE:HG12	5:CE:90:VAL:N	2.25	0.51
6:CF:38:GLU:HG3	6:CF:65:VAL:HA	1.92	0.51
7:CG:92:SER:OG	7:CG:94:ARG:NE	2.43	0.51
9:CI:13:ALA:HA	9:CI:67:GLY:C	2.30	0.51
9:CI:53:VAL:HG12	9:CI:95:LYS:HE3	1.92	0.51
9:CI:28:VAL:HA	9:CI:63:ILE:O	2.11	0.51
10:CJ:64:GLU:N	14:CN:59:ALA:HB2	2.25	0.51
1:CA:692:U:H5	11:CK:26:ASN:ND2	2.08	0.51
12:CL:60:LEU:HD23	12:CL:64:TYR:C	2.31	0.51
1:CA:656:C:H4'	15:CO:62:GLN:NE2	2.25	0.51
16:CP:23:ASP:HB3	16:CP:26:ARG:HG3	1.93	0.51
17:CQ:47:PRO:HG2	17:CQ:48:GLU:OE2	2.11	0.51
28:D2:26:ARG:NH2	54:DX:7:VAL:N	2.56	0.51
31:D5:15:ARG:HA	31:D5:18:ALA:CB	2.40	0.51
34:D8:58:ILE:HG22	34:D8:58:ILE:O	2.11	0.51
35:DA:1166:C:H2'	35:DA:1167:U:H6	1.75	0.51
35:DA:1190:G:C4'	46:DP:35:HIS:HB3	2.40	0.51
35:DA:1227:G:O2'	35:DA:1228:G:H5'	2.10	0.51
35:DA:1352:U:O2'	35:DA:1353:A:H5'	2.11	0.51
35:DA:1608:A:O3'	35:DA:1609:A:H3'	2.10	0.51
35:DA:1708:C:O2'	35:DA:1709:U:H5'	2.09	0.51
35:DA:2314:C:O2'	35:DA:2315:G:H5'	2.10	0.51
35:DA:2553:G:H2'	35:DA:2554:U:H4'	1.91	0.51
35:DA:2491:U:C5'	35:DA:2570:G:H5''	2.22	0.51
35:DA:2821:A:OP2	48:DR:2:ARG:NH1	2.43	0.51
35:DA:685:A:C5	35:DA:774:A:C2	2.98	0.51
35:DA:873:G:H1	35:DA:904:C:H42	1.58	0.51
36:DB:104:U:H4'	56:DZ:72:ARG:HH11	1.74	0.51
37:DC:86:ALA:O	37:DC:94:VAL:HG21	2.11	0.51
39:DE:171:GLU:O	39:DE:173:VAL:HG23	2.10	0.51
40:DF:39:TRP:CE3	40:DF:40:GLN:HG2	2.45	0.51
41:DG:12:TYR:HA	41:DG:16:ARG:HG3	1.92	0.51
43:DI:29:TYR:HD1	43:DI:33:ARG:HE	1.56	0.51
43:DI:88:ILE:HG23	43:DI:89:TYR:H	1.75	0.51
46:DP:57:THR:C	46:DP:59:LEU:H	2.13	0.51
46:DP:71:VAL:HG22	46:DP:72:PRO:CD	2.40	0.51
49:DS:92:TYR:HD1	49:DS:93:LYS:N	2.06	0.51
50:DT:102:ILE:HG12	50:DT:103:ARG:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:DU:88:ILE:HA	51:DU:90:VAL:HG23	1.92	0.51
52:DV:19:LYS:CE	52:DV:19:LYS:HA	2.41	0.51
52:DV:89:GLN:HE21	52:DV:89:GLN:HA	1.74	0.51
31:D5:20:ARG:NH1	53:DW:15:ARG:NH2	2.50	0.51
54:DX:88:LYS:C	54:DX:90:GLU:N	2.63	0.51
28:D2:36:ARG:HH12	54:DX:92:LEU:HD22	1.74	0.51
56:DZ:9:TYR:HE1	56:DZ:61:LEU:HD13	1.76	0.51
1:AA:1116:C:H3'	1:AA:1117:G:H5''	1.90	0.51
1:AA:1178:G:N2	1:AA:1180:A:H3'	2.26	0.51
1:AA:1255:G:H2'	1:AA:1255:G:N3	2.24	0.51
1:AA:223:U:H2'	1:AA:224:C:H6	1.76	0.51
1:AA:509:A:H2'	1:AA:510:A:C8	2.46	0.51
2:AB:69:LEU:HB3	2:AB:162:ILE:CG2	2.40	0.51
4:AD:176:LEU:CG	4:AD:177:ASP:N	2.63	0.51
5:AE:8:GLU:N	5:AE:34:VAL:HG13	2.25	0.51
6:AF:26:ILE:O	6:AF:29:ALA:HB3	2.11	0.51
6:AF:61:LEU:O	6:AF:62:TRP:CB	2.58	0.51
10:AJ:50:ILE:HD11	14:AN:41:ARG:CD	2.35	0.51
11:AK:123:LYS:O	11:AK:124:LYS:C	2.49	0.51
15:AO:70:LEU:HG	15:AO:78:TYR:HB2	1.93	0.51
16:AP:74:LEU:O	16:AP:79:VAL:HG23	2.11	0.51
18:AR:36:ASN:O	18:AR:39:VAL:HG23	2.10	0.51
20:AT:38:LYS:O	20:AT:39:LYS:C	2.49	0.51
20:AT:41:ILE:C	20:AT:43:LEU:N	2.63	0.51
25:AY:20:VAL:O	25:AY:23:HIS:HB3	2.11	0.51
34:B8:34:TRP:CZ3	34:B8:41:ILE:HG23	2.46	0.51
34:B8:58:ILE:HG22	34:B8:58:ILE:O	2.11	0.51
35:BA:1755:A:H2	35:BA:2716:U:C1'	2.20	0.51
35:BA:1829:A:H2'	35:BA:1830:C:O4'	2.10	0.51
1:AA:1418:A:H2	35:BA:1948:G:N3	2.09	0.51
35:BA:2552:U:C2	35:BA:2554:U:H5'	2.46	0.51
35:BA:2839:G:H2'	35:BA:2840:C:C6	2.45	0.51
35:BA:481:G:C2'	35:BA:482:A:OP2	2.59	0.51
35:BA:516:C:O2'	35:BA:517:C:H5'	2.11	0.51
35:BA:644:A:O2'	35:BA:645:C:H5''	2.10	0.51
35:BA:921:G:H2'	35:BA:922:U:C6	2.45	0.51
38:BD:158:ALA:O	38:BD:159:ALA:C	2.49	0.51
38:BD:165:ILE:CD1	38:BD:165:ILE:N	2.70	0.51
38:BD:203:ASN:O	38:BD:204:ILE:O	2.29	0.51
35:BA:1569:A:O2'	38:BD:38:LYS:HE2	2.11	0.51
38:BD:35:LYS:CG	38:BD:64:ILE:H	2.13	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:197:ILE:CG1	39:BE:199:ARG:HH12	2.23	0.51
39:BE:1:MET:HB3	39:BE:200:GLU:OE1	2.10	0.51
39:BE:68:ALA:C	39:BE:70:ALA:H	2.14	0.51
40:BF:114:VAL:O	40:BF:115:ALA:C	2.48	0.51
40:BF:65:TRP:CZ3	40:BF:72:ARG:HB3	2.46	0.51
44:BN:24:GLY:HA2	44:BN:27:ALA:HB3	1.92	0.51
44:BN:34:LEU:CD2	44:BN:120:LEU:HD23	2.41	0.51
44:BN:46:VAL:CG1	44:BN:48:MET:HG3	2.40	0.51
47:BQ:17:LEU:HD21	47:BQ:41:TRP:HE1	1.76	0.51
50:BT:100:TYR:O	50:BT:103:ARG:N	2.43	0.51
50:BT:54:ARG:HA	50:BT:59:THR:OG1	2.09	0.51
35:BA:2875:C:O2'	50:BT:5:ALA:HB3	2.10	0.51
52:BV:4:ILE:CG1	52:BV:40:LEU:HD11	2.40	0.51
54:BX:55:ASN:ND2	54:BX:55:ASN:N	2.57	0.51
54:BX:65:ARG:NE	54:BX:66:LEU:N	2.58	0.51
54:BX:77:LYS:CD	54:BX:78:LYS:H	2.24	0.51
56:BZ:109:ALA:C	56:BZ:110:GLY:O	2.46	0.51
56:BZ:94:GLU:OE2	56:BZ:95:PRO:HG2	2.10	0.51
1:CA:1220:G:O2'	1:CA:1221:G:H5'	2.10	0.51
1:CA:1409:C:H2'	1:CA:1410:G:H8	1.76	0.51
1:CA:1442(A):G:C3'	1:CA:1442(B):A:C5'	2.81	0.51
1:CA:1516:G:H22	1:CA:1518:A:H3'	1.76	0.51
1:CA:551:U:C2	1:CA:552:U:C5	2.99	0.51
1:CA:552:U:H4'	12:CL:87:GLY:H	1.75	0.51
1:CA:707:C:O2'	1:CA:708:C:H5'	2.10	0.51
2:CB:165:VAL:O	2:CB:187:LEU:O	2.28	0.51
2:CB:80:ILE:HD11	2:CB:215:LEU:HD12	1.92	0.51
4:CD:163:GLU:C	4:CD:165:MET:N	2.61	0.51
5:CE:102:ALA:HB1	5:CE:106:PRO:HG2	1.91	0.51
7:CG:21:VAL:HG23	7:CG:22:LEU:N	2.25	0.51
7:CG:47:CYS:O	7:CG:58:PRO:HG3	2.10	0.51
8:CH:11:THR:HG23	8:CH:14:ARG:NH1	2.21	0.51
1:CA:691:G:H1	11:CK:52:GLY:HA2	1.75	0.51
1:CA:880:C:OP2	12:CL:6:THR:HG21	2.11	0.51
15:CO:82:ILE:C	15:CO:82:ILE:HD13	2.31	0.51
25:CY:142:LYS:O	25:CY:145:LYS:HB3	2.10	0.51
27:D1:37:ILE:O	27:D1:38:SER:HB2	2.11	0.51
28:D2:12:GLU:O	28:D2:14:ARG:NH1	2.43	0.51
31:D5:25:LEU:HD12	53:DW:23:LEU:CD1	2.41	0.51
34:D8:53:PRO:C	34:D8:55:ALA:N	2.64	0.51
35:DA:1112:G:C1'	35:DA:1113:U:OP1	2.56	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1272:A:OP2	35:DA:1647:G:OP1	2.28	0.51
35:DA:1568:G:H4'	38:DD:59:LYS:CG	2.41	0.51
35:DA:1788:C:O5'	35:DA:1788:C:H6	1.93	0.51
35:DA:1829:A:H2'	35:DA:1830:C:O4'	2.10	0.51
35:DA:2097:C:O2'	35:DA:2098:U:H5'	2.09	0.51
35:DA:2200:C:N4	35:DA:2223:G:H1	2.07	0.51
35:DA:2441:C:O2	35:DA:2441:C:H2'	2.10	0.51
35:DA:2617:C:O2'	35:DA:2618:G:H5'	2.11	0.51
35:DA:261:G:C2	35:DA:262:A:C8	2.98	0.51
35:DA:2663:G:O2'	35:DA:2664:G:H5'	2.10	0.51
35:DA:453:C:H1'	35:DA:457:A:C2	2.45	0.51
37:DC:89:ALA:HA	37:DC:153:ILE:CB	2.41	0.51
38:DD:201:HIS:C	38:DD:203:ASN:H	2.13	0.51
42:DH:98:LEU:HD12	42:DH:102:ALA:O	2.10	0.51
42:DH:60:ARG:O	42:DH:64:LEU:HG	2.10	0.51
45:DO:20:MET:HE3	45:DO:44:LYS:HG3	1.91	0.51
45:DO:47:ILE:HG23	45:DO:48:PRO:CD	2.39	0.51
45:DO:63:VAL:HG11	45:DO:85:VAL:HG23	1.92	0.51
47:DQ:119:ARG:HG3	47:DQ:119:ARG:HH11	1.76	0.51
50:DT:96:ARG:HB3	50:DT:96:ARG:CZ	2.40	0.51
52:DV:78:LYS:CD	52:DV:79:VAL:N	2.66	0.51
1:AA:1054:C:H2'	1:AA:1054:C:O2	2.11	0.51
1:AA:1054:C:P	1:AA:1197:G:OP2	2.68	0.51
1:AA:1416:G:H2'	1:AA:1417:G:C8	2.45	0.51
1:AA:1440:C:O2'	1:AA:1441:G:H5'	2.10	0.51
1:AA:192:U:H4'	20:AT:102:GLY:O	2.11	0.51
1:AA:324:G:N2	1:AA:327:A:C8	2.79	0.51
1:AA:410:G:N1	1:AA:429:U:O2	2.44	0.51
1:AA:525:C:H2'	1:AA:526:C:H6	1.76	0.51
1:AA:601:C:H2'	1:AA:602:A:C8	2.43	0.51
1:AA:651:C:H2'	1:AA:652:U:C6	2.46	0.51
1:AA:707:C:O2'	1:AA:708:C:H5'	2.11	0.51
1:AA:908:A:C4	1:AA:909:A:N7	2.79	0.51
1:AA:950:U:H1'	1:AA:971:G:N7	2.26	0.51
2:AB:167:PRO:O	2:AB:168:THR:C	2.48	0.51
6:AF:83:ASP:O	6:AF:85:VAL:N	2.43	0.51
8:AH:114:THR:HG23	8:AH:116:LYS:O	2.11	0.51
8:AH:9:MET:O	8:AH:12:ARG:N	2.42	0.51
12:AL:32:PHE:HE1	12:AL:86:ARG:HG3	1.76	0.51
13:AM:74:VAL:HA	13:AM:77:ASN:HD22	1.75	0.51
16:AP:9:PHE:HE2	16:AP:18:ARG:NE	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AP:19:ILE:CG2	16:AP:36:ILE:HG13	2.39	0.51
17:AQ:99:SER:C	17:AQ:100:LYS:HE2	2.31	0.51
17:AQ:29:HIS:HE1	17:AQ:31:LEU:HB3	1.74	0.51
17:AQ:31:LEU:HG	17:AQ:32:TYR:CD2	2.46	0.51
25:AY:3:LEU:HB3	25:AY:7:TYR:CE1	2.45	0.51
25:AY:64:ARG:N	25:AY:64:ARG:CZ	2.74	0.51
27:B1:10:LYS:HG3	27:B1:11:ARG:N	2.25	0.51
27:B1:87:PRO:HA	27:B1:90:ILE:HG12	1.93	0.51
34:B8:31:HIS:O	34:B8:32:LEU:C	2.48	0.51
35:BA:1338:G:H5'	35:BA:1339:G:OP2	2.10	0.51
35:BA:1367:A:H3'	35:BA:1368:G:O4'	2.11	0.51
35:BA:1486:A:N6	35:BA:1504:C:H42	2.09	0.51
35:BA:2119:A:C2'	35:BA:2120:G:H5''	2.40	0.51
35:BA:2590:A:O3'	38:BD:239:ARG:HG3	2.10	0.51
35:BA:2692:C:O2'	35:BA:2693:A:H5'	2.11	0.51
35:BA:272(C):G:H2'	35:BA:272(D):G:C8	2.40	0.51
35:BA:1050:A:C4	35:BA:2751:G:C2	2.99	0.51
35:BA:2842:G:O2'	35:BA:2843:G:H5'	2.11	0.51
35:BA:528:A:OP2	44:BN:114:ARG:HD2	2.10	0.51
35:BA:838:C:O2'	35:BA:839:U:H5'	2.11	0.51
38:BD:173:VAL:HG12	38:BD:185:VAL:O	2.11	0.51
39:BE:103:ASP:OD1	39:BE:168:MET:HG2	2.11	0.51
39:BE:8:LYS:HD2	39:BE:189:PRO:O	2.10	0.51
39:BE:47:VAL:HG23	39:BE:84:PHE:O	2.09	0.51
40:BF:43:LYS:HA	40:BF:98:SER:HA	1.92	0.51
41:BG:57:ALA:O	41:BG:60:LEU:HB3	2.10	0.51
41:BG:93:THR:CG2	41:BG:95:ARG:HG3	2.39	0.51
42:BH:85:LYS:O	42:BH:132:ARG:CA	2.56	0.51
43:BI:92:VAL:HG12	43:BI:119:PRO:O	2.11	0.51
49:BS:29:PHE:CD1	49:BS:29:PHE:C	2.84	0.51
50:BT:28:VAL:HG11	50:BT:46:GLU:OE1	2.11	0.51
50:BT:96:ARG:CG	50:BT:96:ARG:NH1	2.71	0.51
52:BV:61:VAL:HG23	52:BV:100:ARG:HG2	1.91	0.51
53:BW:17:VAL:O	53:BW:18:ARG:C	2.49	0.51
54:BX:7:VAL:HA	54:BX:31:HIS:HB2	1.92	0.51
55:BY:37:VAL:O	55:BY:38:ILE:CB	2.59	0.51
1:CA:1131:G:H2'	1:CA:1132:C:C5	2.45	0.51
1:CA:1253:G:O2'	1:CA:1254:C:H5'	2.10	0.51
1:CA:1307:U:H2'	1:CA:1308:U:H6	1.71	0.51
1:CA:59:A:C5'	1:CA:60:A:C5'	2.86	0.51
1:CA:874:G:H2'	1:CA:875:C:H6	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:30:ARG:HD2	14:CN:38:GLY:HA3	1.93	0.51
4:CD:68:TYR:CZ	4:CD:97:LEU:HD22	2.46	0.51
5:CE:146:ALA:C	5:CE:148:VAL:N	2.61	0.51
5:CE:72:GLN:HE22	5:CE:77:PRO:CD	2.22	0.51
6:CF:67:MET:CE	6:CF:72:VAL:H	2.24	0.51
7:CG:15:ASP:OD2	7:CG:16:LEU:N	2.43	0.51
7:CG:88:PRO:O	7:CG:89:MET:HB3	2.09	0.51
8:CH:112:LEU:C	8:CH:112:LEU:HD12	2.30	0.51
8:CH:69:ARG:HB2	8:CH:74:PRO:HA	1.93	0.51
9:CI:4:TYR:HD1	9:CI:4:TYR:N	2.08	0.51
10:CJ:16:LEU:HA	10:CJ:19:SER:OG	2.10	0.51
16:CP:72:ARG:O	16:CP:74:LEU:N	2.43	0.51
1:CA:1325:C:H4'	21:CU:17:THR:HG21	1.93	0.51
22:CV:31:U:H2'	22:CV:31:U:O2	2.09	0.51
30:D4:43:TYR:C	30:D4:45:GLY:H	2.14	0.51
34:D8:48:PHE:H	34:D8:48:PHE:HD1	1.54	0.51
35:DA:1015:G:H2'	35:DA:1015:G:N3	2.24	0.51
35:DA:1131:G:H1'	35:DA:1132:A:C8	2.46	0.51
35:DA:1410:G:C6	35:DA:1411:C:N4	2.79	0.51
35:DA:1604:C:O2'	35:DA:1605:C:H5'	2.10	0.51
35:DA:179:G:O2'	35:DA:180:G:H5'	2.11	0.51
35:DA:2187:G:O2'	35:DA:2188:C:H5'	2.11	0.51
35:DA:323:G:O2'	35:DA:1205:U:N3	2.44	0.51
38:DD:132:PRO:O	38:DD:136:ILE:HG13	2.10	0.51
38:DD:213:ARG:O	38:DD:215:LEU:N	2.43	0.51
41:DG:106:LEU:O	41:DG:108:ASN:N	2.44	0.51
44:DN:126:PRO:O	44:DN:127:ASP:CB	2.58	0.51
45:DO:1:MET:CE	45:DO:1:MET:H3	2.21	0.51
45:DO:44:LYS:O	45:DO:45:GLU:HB3	2.09	0.51
46:DP:90:ARG:HH11	46:DP:91:PHE:HB3	1.75	0.51
47:DQ:20:ALA:CB	47:DQ:98:LYS:HB3	2.40	0.51
49:DS:29:PHE:HB3	49:DS:36:TYR:HB2	1.92	0.51
50:DT:27:THR:HG1	50:DT:87:ASP:HA	1.74	0.51
54:DX:77:LYS:CD	54:DX:78:LYS:H	2.23	0.51
56:DZ:141:VAL:O	56:DZ:144:LEU:HD23	2.11	0.51
56:DZ:96:VAL:CG2	56:DZ:97:GLU:N	2.74	0.51
1:AA:1046:A:H3'	1:AA:1047:G:C8	2.44	0.51
1:AA:1261:A:H62	1:AA:1274:G:N2	2.09	0.51
1:AA:1325:C:H4'	21:AU:17:THR:HG21	1.93	0.51
1:AA:1478:C:H2'	1:AA:1479:C:H6	1.72	0.51
1:AA:185:A:H61	1:AA:192:U:H3	1.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:406:G:C2	1:AA:407:G:C8	2.99	0.51
1:AA:70:G:H2'	1:AA:71:C:C6	2.45	0.51
4:AD:192:GLU:N	4:AD:192:GLU:OE2	2.44	0.51
4:AD:25:ARG:C	4:AD:27:TYR:H	2.13	0.51
5:AE:146:ALA:C	5:AE:148:VAL:N	2.64	0.51
5:AE:26:PHE:O	5:AE:27:ARG:HB2	2.11	0.51
8:AH:40:ALA:HB2	8:AH:45:ILE:CD1	2.41	0.51
9:AI:82:ALA:O	9:AI:96:LEU:HD21	2.11	0.51
10:AJ:45:ARG:HH11	10:AJ:45:ARG:HG3	1.75	0.51
11:AK:20:TYR:C	11:AK:21:ILE:HD12	2.31	0.51
12:AL:117:ARG:O	12:AL:119:LYS:O	2.29	0.51
15:AO:17:ARG:NH1	15:AO:77:ARG:NH1	2.58	0.51
16:AP:82:GLN:NE2	16:AP:82:GLN:N	2.44	0.51
18:AR:44:LEU:O	18:AR:45:SER:CB	2.59	0.51
19:AS:16:LEU:N	19:AS:16:LEU:CD1	2.74	0.51
25:AY:137:LEU:O	25:AY:140:LEU:HB3	2.11	0.51
25:AY:73:GLN:O	25:AY:77:LYS:HG2	2.10	0.51
27:B1:11:ARG:NE	27:B1:61:ARG:H	2.09	0.51
27:B1:20:ARG:CD	27:B1:20:ARG:N	2.71	0.51
27:B1:76:ARG:HA	27:B1:78:LYS:HZ3	1.75	0.51
27:B1:83:GLU:HG2	27:B1:86:SER:H	1.76	0.51
35:BA:1245:G:H3'	46:BP:16:ARG:NH2	2.22	0.51
35:BA:1378:A:C4'	35:BA:1379:A:OP1	2.55	0.51
35:BA:1612:C:C2'	35:BA:1613:G:O5'	2.59	0.51
35:BA:154(A):C:C5	35:BA:171:G:N1	2.78	0.51
35:BA:17:G:O2'	51:BU:25:TRP:HZ3	1.93	0.51
35:BA:1904:G:N2	35:BA:1905:C:H1'	2.25	0.51
35:BA:2128:C:H5'	35:BA:2173:A:C2	2.45	0.51
35:BA:2713:A:C2'	35:BA:2714:G:H5'	2.41	0.51
35:BA:305:U:O2'	35:BA:306:U:H5'	2.11	0.51
35:BA:460:A:C6	35:BA:470:A:C8	2.98	0.51
35:BA:785:G:H2'	35:BA:786:C:H6	1.75	0.51
35:BA:876:C:H2'	35:BA:877:U:C6	2.46	0.51
38:BD:24:ILE:O	38:BD:26:LYS:N	2.43	0.51
39:BE:142:GLY:C	39:BE:143:ASN:HD22	2.14	0.51
39:BE:47:VAL:HG12	39:BE:49:LEU:CD2	2.40	0.51
41:BG:13:GLU:O	41:BG:14:GLU:HB3	2.10	0.51
43:BI:48:GLU:O	43:BI:51:ILE:HB	2.10	0.51
47:BQ:110:THR:OG1	47:BQ:113:GLN:HG3	2.10	0.51
47:BQ:119:ARG:HG3	47:BQ:119:ARG:HH11	1.76	0.51
48:BR:17:ARG:CG	48:BR:17:ARG:HH11	2.17	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BS:65:VAL:CG1	49:BS:69:VAL:HB	2.41	0.51
51:BU:98:LEU:O	51:BU:101:ARG:O	2.28	0.51
56:BZ:129:SER:O	56:BZ:131:ARG:N	2.42	0.51
1:CA:1277:C:O2'	1:CA:1278:U:H5'	2.10	0.51
1:CA:1430:C:H2'	1:CA:1431:C:C6	2.45	0.51
1:CA:199:G:H2'	1:CA:200:G:C8	2.46	0.51
1:CA:19:C:P	5:CE:127:ASN:HD22	2.34	0.51
1:CA:309:G:H2'	1:CA:310:G:H8	1.76	0.51
1:CA:309:G:O2'	1:CA:310:G:H5'	2.10	0.51
1:CA:774:G:OP1	38:DD:202:LYS:NZ	2.40	0.51
1:CA:859:A:O2'	1:CA:860:A:H5'	2.10	0.51
1:CA:914:A:O2'	1:CA:915:A:H5'	2.11	0.51
1:CA:985:C:H2'	1:CA:986:A:H8	1.74	0.51
2:CB:58:ILE:O	2:CB:61:LEU:HB3	2.11	0.51
3:CC:106:VAL:HG12	3:CC:108:ASN:H	1.75	0.51
4:CD:13:ARG:CG	4:CD:14:ARG:H	2.24	0.51
6:CF:61:LEU:O	6:CF:62:TRP:CB	2.59	0.51
6:CF:66:GLU:O	6:CF:67:MET:HB3	2.11	0.51
8:CH:107:LEU:HD23	8:CH:107:LEU:C	2.31	0.51
8:CH:10:LEU:H	8:CH:10:LEU:CD2	2.18	0.51
14:CN:4:LYS:HA	14:CN:7:ILE:CD1	2.40	0.51
14:CN:8:GLU:C	14:CN:12:ARG:HD3	2.31	0.51
15:CO:78:TYR:C	15:CO:80:ALA:N	2.64	0.51
1:CA:378:G:P	16:CP:3:LYS:HZ1	2.33	0.51
24:CX:19:U:C2'	24:CX:19:U:O2	2.59	0.51
25:CY:25:LEU:O	25:CY:179:LYS:HE2	2.11	0.51
27:D1:10:LYS:O	27:D1:13:ILE:CG2	2.53	0.51
27:D1:25:LYS:CG	27:D1:37:ILE:HG21	2.32	0.51
28:D2:53:LEU:HB2	35:DA:76:C:O3'	2.10	0.51
29:D3:4:LEU:HD12	29:D3:39:ASP:OD1	2.10	0.51
31:D5:16:ARG:NH1	31:D5:16:ARG:CG	2.74	0.51
35:DA:104:U:H6	35:DA:104:U:O5'	1.94	0.51
35:DA:1131:G:H1'	35:DA:1132:A:H8	1.75	0.51
35:DA:1686:C:H42	35:DA:1702:G:H1	1.59	0.51
35:DA:1766:U:O2'	35:DA:1767:C:H5'	2.11	0.51
35:DA:2315:G:H2'	35:DA:2316:C:H6	1.69	0.51
35:DA:272(C):G:O2'	35:DA:272(D):G:H5'	2.10	0.51
35:DA:863:A:H4'	36:DB:101:G:N2	2.25	0.51
35:DA:871:U:H4'	47:DQ:69:PHE:CE2	2.46	0.51
35:DA:911:A:C4	47:DQ:9:TYR:OH	2.61	0.51
37:DC:214:VAL:C	37:DC:216:THR:N	2.63	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DC:49:ILE:C	37:DC:51:PRO:HD3	2.31	0.51
38:DD:125:ILE:H	38:DD:125:ILE:HD12	1.76	0.51
39:DE:4:ILE:HD12	39:DE:31:CYS:SG	2.51	0.51
36:DB:41:U:N3	41:DG:70:VAL:O	2.43	0.51
42:DH:12:PRO:N	42:DH:15:VAL:HG21	2.26	0.51
35:DA:598:G:C5'	46:DP:15:ARG:HD2	2.37	0.51
46:DP:79:ARG:O	46:DP:111:ARG:HB2	2.10	0.51
47:DQ:81:VAL:CG2	47:DQ:82:ARG:HG2	2.41	0.51
49:DS:99:LYS:O	49:DS:100:ALA:C	2.49	0.51
51:DU:90:VAL:HG13	52:DV:39:LEU:CB	2.36	0.51
52:DV:72:VAL:HG13	52:DV:88:ARG:HH22	1.75	0.51
35:DA:2012:G:O3'	53:DW:96:ILE:HD11	2.10	0.51
54:DX:31:HIS:CG	54:DX:32:PRO:HD2	2.46	0.51
54:DX:35:THR:HG23	54:DX:36:LYS:H	1.74	0.51
54:DX:52:VAL:O	54:DX:53:LYS:HB3	2.11	0.51
55:DY:83:THR:HG22	55:DY:84:ARG:H	1.76	0.51
56:DZ:144:LEU:CD1	56:DZ:149:SER:HA	2.40	0.51
1:AA:328:C:O2'	1:AA:329:A:OP2	2.28	0.51
1:AA:645:C:H2'	1:AA:646:U:H6	1.76	0.51
2:AB:111:ARG:O	2:AB:115:LEU:N	2.44	0.51
2:AB:234:PRO:O	2:AB:236:TYR:N	2.44	0.51
2:AB:70:PHE:O	2:AB:92:TYR:HA	2.11	0.51
3:AC:180:ALA:O	3:AC:205:GLY:O	2.28	0.51
4:AD:119:GLN:HB3	4:AD:120:LEU:HD12	1.91	0.51
5:AE:129:ILE:O	5:AE:130:ASN:C	2.49	0.51
8:AH:114:THR:HG22	8:AH:117:GLY:O	2.11	0.51
11:AK:125:PHE:N	11:AK:125:PHE:HD1	2.08	0.51
11:AK:21:ILE:HD13	11:AK:84:VAL:CG1	2.38	0.51
18:AR:65:ILE:HD12	18:AR:66:LEU:H	1.75	0.51
13:AM:91:ARG:NH1	19:AS:81:ARG:HH22	1.90	0.51
20:AT:90:GLN:C	20:AT:93:GLU:OE2	2.49	0.51
25:AY:143:LEU:HD23	25:AY:143:LEU:C	2.31	0.51
26:B0:23:VAL:HG12	26:B0:24:LYS:N	2.24	0.51
29:B3:45:GLY:HA3	35:BA:852:G:H5'	1.93	0.51
32:B6:47:THR:HG22	32:B6:48:VAL:N	2.25	0.51
35:BA:1259:G:H2'	35:BA:1260:G:H8	1.75	0.51
35:BA:1268:A:C6	35:BA:2013:A:C8	2.99	0.51
35:BA:132:G:O2'	35:BA:133:C:H5'	2.11	0.51
35:BA:1387:C:H4'	35:BA:1469:A:O4'	2.11	0.51
35:BA:1914:C:H3'	35:BA:1914:C:O2	2.10	0.51
35:BA:2039:C:C2	35:BA:2040:C:C5	2.99	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2251:G:N3	35:BA:2450:A:H1'	2.25	0.51
35:BA:2393:A:O2'	35:BA:2394:C:H5'	2.10	0.51
35:BA:2431:U:H2'	35:BA:2433:A:OP2	2.10	0.51
35:BA:2892:A:C4	35:BA:2893:G:H1'	2.46	0.51
35:BA:563:G:C6	35:BA:564:C:N4	2.79	0.51
35:BA:587:C:H3'	46:BP:33:ARG:NH1	2.20	0.51
35:BA:61:G:C2'	35:BA:62:C:H5'	2.41	0.51
35:BA:950:G:O2'	35:BA:951:C:H5'	2.11	0.51
38:BD:35:LYS:HG2	38:BD:64:ILE:HG22	1.92	0.51
38:BD:65:ILE:C	38:BD:65:ILE:CD1	2.77	0.51
39:BE:5:LEU:CD2	39:BE:197:ILE:HG22	2.41	0.51
40:BF:155:LEU:CD1	40:BF:174:VAL:HB	2.40	0.51
41:BG:16:ARG:NH1	41:BG:16:ARG:HG3	2.26	0.51
41:BG:57:ALA:HB2	41:BG:90:LEU:CD2	2.40	0.51
42:BH:97:ARG:O	42:BH:125:VAL:HG21	2.11	0.51
43:BI:91:SER:H	43:BI:121:LYS:HE3	1.75	0.51
43:BI:94:ALA:HA	43:BI:97:ILE:CB	2.41	0.51
44:BN:26:LEU:CD1	44:BN:30:ILE:HD11	2.41	0.51
47:BQ:74:TYR:HB3	47:BQ:91:GLU:CD	2.32	0.51
50:BT:108:ARG:O	50:BT:112:ARG:HG3	2.09	0.51
50:BT:57:PHE:O	50:BT:59:THR:HG22	2.10	0.51
45:BO:76:ALA:HB3	50:BT:75:ILE:HD13	1.92	0.51
52:BV:32:THR:CG2	52:BV:33:VAL:N	2.73	0.51
53:BW:80:PRO:HD2	53:BW:100:THR:HG21	1.93	0.51
31:B5:25:LEU:HD12	53:BW:23:LEU:CD1	2.41	0.51
35:BA:1615:C:O2	53:BW:87:PRO:HG2	2.10	0.51
54:BX:88:LYS:C	54:BX:90:GLU:N	2.63	0.51
56:BZ:52:SER:HB3	56:BZ:54:HIS:CD2	2.45	0.51
1:CA:1166:G:H2'	1:CA:1169:A:OP2	2.11	0.51
1:CA:1264:C:H2'	1:CA:1265:G:C8	2.39	0.51
1:CA:1256:A:N6	1:CA:1278:U:H1'	2.26	0.51
1:CA:556:C:C2'	1:CA:557:G:H5'	2.41	0.51
1:CA:603:U:H2'	1:CA:604:G:H8	1.76	0.51
1:CA:627:G:O2'	1:CA:628:G:H5'	2.10	0.51
1:CA:705:U:C5	1:CA:706:A:C5	2.98	0.51
2:CB:115:LEU:HD21	2:CB:153:ARG:HE	1.76	0.51
3:CC:186:PHE:HD1	3:CC:198:VAL:O	1.93	0.51
3:CC:60:ALA:CB	3:CC:63:ASN:HD21	2.24	0.51
3:CC:8:ILE:O	3:CC:10:PHE:N	2.44	0.51
4:CD:80:GLU:C	4:CD:84:LYS:HZ3	2.14	0.51
5:CE:79:GLU:OE1	5:CE:79:GLU:N	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:939:G:H5''	7:CG:102:ARG:NH1	2.25	0.51
8:CH:119:LEU:HD12	8:CH:123:GLU:C	2.31	0.51
15:CO:43:LEU:C	15:CO:45:VAL:H	2.14	0.51
15:CO:87:ILE:HG22	15:CO:88:ARG:N	2.26	0.51
1:CA:626:U:H5''	16:CP:38:TYR:CD2	2.46	0.51
18:CR:29:PHE:HD1	18:CR:39:VAL:HG11	1.76	0.51
18:CR:44:LEU:O	18:CR:45:SER:CB	2.58	0.51
25:CY:116:ARG:O	25:CY:117:ALA:C	2.49	0.51
26:D0:72:ARG:HB3	26:D0:75:LEU:HB3	1.91	0.51
27:D1:87:PRO:C	27:D1:89:GLU:OE2	2.50	0.51
27:D1:87:PRO:C	27:D1:90:ILE:HG12	2.30	0.51
28:D2:12:GLU:O	28:D2:12:GLU:CD	2.49	0.51
33:D7:29:LYS:NZ	33:D7:32:LYS:NZ	2.59	0.51
35:DA:1202:C:C2'	35:DA:1203:G:H5'	2.41	0.51
35:DA:1387:C:H4'	35:DA:1469:A:O4'	2.11	0.51
35:DA:1485:G:H2'	35:DA:1486:A:C8	2.45	0.51
35:DA:1707:G:H1'	35:DA:1756:G:N3	2.25	0.51
35:DA:201:C:C2'	35:DA:202:U:H5'	2.39	0.51
35:DA:2699:C:O2'	35:DA:2700:C:H5'	2.11	0.51
35:DA:2712(A):A:OP2	35:DA:2714:G:OP2	2.28	0.51
35:DA:2821:A:P	48:DR:2:ARG:NH2	2.63	0.51
35:DA:2850:A:H5'	35:DA:2868:A:H2	1.75	0.51
35:DA:2889:C:H2'	35:DA:2889:C:O2	2.09	0.51
35:DA:494:G:O2'	35:DA:495:G:H5'	2.11	0.51
35:DA:587:C:O2'	35:DA:588:U:OP2	2.24	0.51
35:DA:675:A:C8	35:DA:804:A:C6	2.98	0.51
36:DB:15:A:C3'	36:DB:16:G:H5'	2.40	0.51
37:DC:214:VAL:O	37:DC:216:THR:N	2.44	0.51
38:DD:127:VAL:HA	38:DD:193:VAL:HG13	1.93	0.51
39:DE:110:GLY:O	48:DR:2:ARG:NH2	2.44	0.51
39:DE:73:GLU:N	39:DE:73:GLU:OE2	2.36	0.51
40:DF:62:ARG:NH2	40:DF:64:ILE:HA	2.26	0.51
41:DG:142:PRO:C	41:DG:144:ILE:H	2.14	0.51
42:DH:88:LEU:O	42:DH:89:ILE:CG2	2.59	0.51
43:DI:127:VAL:O	43:DI:127:VAL:HG12	2.11	0.51
43:DI:29:TYR:CE1	43:DI:33:ARG:NE	2.79	0.51
44:DN:62:VAL:CG2	44:DN:66:LYS:HB2	2.41	0.51
44:DN:70:LYS:HB3	44:DN:87:LEU:HB2	1.91	0.51
46:DP:106:LEU:C	46:DP:107:LYS:HG2	2.31	0.51
46:DP:80:TYR:CZ	46:DP:111:ARG:HG2	2.45	0.51
54:DX:33:LYS:O	54:DX:35:THR:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:DX:55:ASN:HB2	54:DX:77:LYS:CD	2.41	0.51
56:DZ:27:VAL:CG1	56:DZ:85:HIS:HE2	2.23	0.51
1:AA:1332:A:H2'	1:AA:1333:A:C8	2.45	0.51
1:AA:355:C:C2	1:AA:356:A:C8	2.99	0.51
1:AA:666:G:H2'	1:AA:667:G:H8	1.76	0.51
1:AA:740:U:O2'	1:AA:741:G:H5'	2.11	0.51
1:AA:751:U:C2'	1:AA:752:G:H5'	2.40	0.51
2:AB:77:ALA:O	2:AB:80:ILE:HG23	2.11	0.51
3:AC:18:TRP:HE3	3:AC:18:TRP:H	1.54	0.51
4:AD:110:PHE:N	4:AD:110:PHE:HD1	2.09	0.51
4:AD:121:VAL:HA	4:AD:126:ILE:HD12	1.91	0.51
5:AE:93:PRO:HA	5:AE:118:ILE:HD12	1.92	0.51
14:AN:33:VAL:HG12	14:AN:34:TYR:N	2.26	0.51
16:AP:18:ARG:O	16:AP:20:VAL:HG12	2.11	0.51
19:AS:72:GLY:C	19:AS:74:PHE:H	2.14	0.51
25:AY:144:ALA:C	25:AY:147:LEU:O	2.49	0.51
25:AY:61:PRO:HG3	25:AY:67:VAL:HG13	1.92	0.51
27:B1:47:GLN:NE2	27:B1:64:ALA:HB2	2.25	0.51
29:B3:3:ARG:HG2	29:B3:38:GLU:OE2	2.11	0.51
35:BA:1141:U:H5''	35:BA:1142(A):A:O4'	2.11	0.51
35:BA:1250:G:H3'	35:BA:1251:C:H5'	1.92	0.51
35:BA:1532:C:O2	35:BA:1532:C:H2'	2.10	0.51
35:BA:1643:G:H2'	35:BA:1644:C:C6	2.46	0.51
35:BA:179:G:O2'	35:BA:180:G:H5'	2.11	0.51
35:BA:1947:C:O2'	35:BA:1948:G:H5'	2.10	0.51
35:BA:2415:G:O3'	46:BP:66:GLY:CA	2.49	0.51
35:BA:2579:C:H4'	39:BE:134:ILE:HG13	1.92	0.51
35:BA:272(C):G:O2'	35:BA:272(D):G:H5'	2.11	0.51
35:BA:862:G:H2'	35:BA:863:A:O4'	2.11	0.51
35:BA:926:A:H2'	35:BA:927:G:C8	2.43	0.51
37:BC:40:THR:HG21	37:BC:215:THR:CB	2.40	0.51
37:BC:86:ALA:O	37:BC:94:VAL:HG21	2.10	0.51
38:BD:21:PHE:O	38:BD:23:GLU:N	2.44	0.51
38:BD:271:ILE:N	38:BD:271:ILE:HD12	2.25	0.51
40:BF:125:LEU:HD12	40:BF:196:LEU:CD2	2.41	0.51
41:BG:31:VAL:CG2	41:BG:32:PRO:HD2	2.41	0.51
42:BH:19:VAL:CG2	42:BH:44:VAL:HA	2.33	0.51
43:BI:111:PRO:HA	43:BI:114:LEU:HD11	1.93	0.51
43:BI:87:LYS:HZ3	43:BI:121:LYS:HG2	1.76	0.51
46:BP:40:SER:C	46:BP:41:ARG:HD2	2.32	0.51
35:BA:631:A:C4'	46:BP:65:ARG:HA	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BR:34:ILE:HG22	48:BR:35:THR:N	2.26	0.51
50:BT:26:ASP:O	50:BT:88:ILE:HB	2.11	0.51
51:BU:47:TYR:HA	51:BU:50:ARG:NH2	2.26	0.51
52:BV:38:LEU:O	52:BV:39:LEU:HD22	2.11	0.51
53:BW:17:VAL:O	53:BW:20:VAL:HG23	2.10	0.51
53:BW:51:LEU:O	53:BW:51:LEU:HD22	2.11	0.51
54:BX:25:LYS:HZ1	54:BX:87:GLN:H	1.56	0.51
55:BY:28:LYS:NZ	55:BY:37:VAL:HG12	2.24	0.51
56:BZ:44:PHE:HE1	56:BZ:48:PHE:CG	2.28	0.51
1:CA:922:G:O2'	1:CA:1398:A:N1	2.42	0.51
1:CA:1404:C:H2'	1:CA:1405:G:H8	1.73	0.51
1:CA:428:G:O4'	1:CA:430:A:C8	2.64	0.51
1:CA:594:G:O2'	1:CA:595:G:H5'	2.11	0.51
1:CA:693:G:H2'	1:CA:694:A:O4'	2.11	0.51
1:CA:777:A:H2'	1:CA:778:G:C8	2.45	0.51
2:CB:12:GLU:C	2:CB:14:GLY:N	2.61	0.51
3:CC:172:ARG:HH11	3:CC:172:ARG:HB3	1.76	0.51
3:CC:35:GLU:HA	3:CC:38:ARG:HG2	1.93	0.51
4:CD:64:LEU:HD23	4:CD:203:VAL:HG21	1.93	0.51
5:CE:146:ALA:O	5:CE:148:VAL:N	2.44	0.51
6:CF:7:ASN:C	6:CF:8:ILE:HG13	2.31	0.51
7:CG:105:VAL:HG12	7:CG:109:ASN:HD21	1.73	0.51
8:CH:40:ALA:HB2	8:CH:45:ILE:HD11	1.92	0.51
8:CH:28:ALA:CB	8:CH:57:PRO:O	2.59	0.51
11:CK:123:LYS:O	11:CK:124:LYS:C	2.50	0.51
13:CM:48:LEU:HD11	13:CM:53:VAL:HG22	1.93	0.51
18:CR:40:LEU:O	18:CR:42:ARG:N	2.44	0.51
18:CR:74:ARG:HH11	18:CR:74:ARG:HG3	1.76	0.51
25:CY:124:GLU:O	25:CY:125:GLY:C	2.50	0.51
25:CY:12:SER:O	25:CY:13:HIS:C	2.49	0.51
25:CY:30:THR:OG1	25:CY:34:ASN:HB2	2.10	0.51
25:CY:55:ILE:CD1	25:CY:55:ILE:N	2.74	0.51
27:D1:11:ARG:HG3	27:D1:61:ARG:O	2.11	0.51
35:DA:1186:G:C2'	35:DA:1187:G:H5'	2.40	0.51
35:DA:1298:C:H3'	35:DA:1299:G:H8	1.76	0.51
35:DA:1496:A:H8	35:DA:1577:C:HO2'	1.57	0.51
35:DA:1688:U:O2	35:DA:1688:U:H2'	2.11	0.51
35:DA:2018:G:C6	35:DA:2019:A:C5	2.99	0.51
35:DA:570:G:H2'	35:DA:2030:A:C5	2.46	0.51
35:DA:203:C:H2'	35:DA:204:A:C8	2.46	0.51
35:DA:2090:G:H2'	35:DA:2091:U:O4'	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2175:C:C3'	35:DA:2176:A:H5''	2.41	0.51
35:DA:256:A:H2'	35:DA:257:A:C8	2.46	0.51
35:DA:2598:A:H5''	38:DD:236:GLY:CA	2.41	0.51
35:DA:2678:C:O2'	35:DA:2679:A:H5'	2.11	0.51
35:DA:2696:U:H2'	35:DA:2697:G:H8	1.74	0.51
35:DA:271(L):U:H4'	35:DA:271(M):G:N7	2.26	0.51
35:DA:2735:G:H2'	35:DA:2736:G:H8	1.76	0.51
35:DA:2870:C:H2'	35:DA:2871:C:H5'	1.93	0.51
35:DA:483:A:H3'	35:DA:484:C:H6	1.76	0.51
35:DA:534:U:H2'	35:DA:535:C:C6	2.46	0.51
35:DA:848:G:C2	35:DA:933:A:H1'	2.46	0.51
39:DE:199:ARG:HB3	39:DE:200:GLU:OE2	2.11	0.51
40:DF:25:PRO:HG3	40:DF:119:ARG:HA	1.91	0.51
40:DF:34:TRP:O	40:DF:37:VAL:N	2.43	0.51
40:DF:7:TYR:HB3	40:DF:16:GLY:C	2.31	0.51
41:DG:101:ILE:HG23	41:DG:102:PHE:N	2.26	0.51
41:DG:27:ASN:O	41:DG:30:GLU:HG2	2.11	0.51
41:DG:45:GLU:C	41:DG:47:LYS:N	2.64	0.51
41:DG:9:ARG:HG2	41:DG:13:GLU:OE1	2.11	0.51
42:DH:84:SER:O	42:DH:85:LYS:CB	2.59	0.51
43:DI:92:VAL:CG1	43:DI:120:ILE:HD12	2.36	0.51
43:DI:71:ILE:HG13	43:DI:72:LEU:HD22	1.92	0.51
43:DI:92:VAL:O	43:DI:92:VAL:HG13	2.11	0.51
45:DO:40:VAL:HA	45:DO:58:VAL:O	2.11	0.51
46:DP:85:LEU:HD23	46:DP:85:LEU:N	2.12	0.51
48:DR:55:ALA:CB	48:DR:79:LEU:HD12	2.41	0.51
49:DS:92:TYR:HB3	49:DS:97:ARG:HH11	1.75	0.51
50:DT:109:GLU:CB	50:DT:113:LYS:HE3	2.24	0.51
50:DT:66:VAL:HG13	50:DT:71:GLY:N	2.26	0.51
52:DV:22:VAL:HG21	52:DV:96:ILE:HB	1.93	0.51
52:DV:39:LEU:HD11	52:DV:53:GLU:CA	2.41	0.51
52:DV:38:LEU:O	52:DV:53:GLU:O	2.28	0.51
56:DZ:122:ARG:HG3	56:DZ:123:ASP:OD2	2.11	0.51
56:DZ:48:PHE:CZ	56:DZ:52:SER:HA	2.45	0.51
56:DZ:48:PHE:O	56:DZ:52:SER:N	2.44	0.51
1:AA:1146:A:H2'	1:AA:1147:C:O5'	2.11	0.51
1:AA:1268:A:H4'	21:AU:20:LYS:N	2.25	0.51
1:AA:1420:C:H3'	1:AA:1420:C:H6	1.76	0.51
1:AA:1433:A:N7	1:AA:1468:A:C6	2.78	0.51
1:AA:1479:C:H2'	1:AA:1480:G:O4'	2.10	0.51
1:AA:18:C:H4'	1:AA:1078:U:O2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:285:G:O2'	1:AA:286:G:H5'	2.11	0.51
1:AA:376:G:H5''	16:AP:5:ARG:CD	2.41	0.51
1:AA:503:C:O2'	1:AA:504:C:H5'	2.11	0.51
1:AA:522:C:N4	1:AA:528:C:N4	2.57	0.51
1:AA:542:G:C4	1:AA:543:C:C5	2.99	0.51
1:AA:66:G:N2	1:AA:172:A:H2	2.08	0.51
2:AB:92:TYR:CD2	2:AB:151:GLY:HA3	2.45	0.51
2:AB:28:PHE:O	2:AB:28:PHE:HD1	1.92	0.51
3:AC:91:LEU:HB2	3:AC:99:VAL:HG21	1.93	0.51
5:AE:149:GLU:C	5:AE:151:LEU:H	2.12	0.51
5:AE:95:ALA:O	5:AE:96:PRO:C	2.50	0.51
7:AG:92:SER:OG	7:AG:93:PRO:HD2	2.11	0.51
11:AK:62:GLN:C	11:AK:64:ALA:H	2.14	0.51
1:AA:1216:G:O3'	14:AN:5:ALA:HB1	2.11	0.51
15:AO:82:ILE:HG12	15:AO:87:ILE:HB	1.92	0.51
17:AQ:92:ARG:O	17:AQ:94:ASN:N	2.44	0.51
18:AR:74:ARG:HE	18:AR:81:PHE:CA	2.21	0.51
27:B1:10:LYS:HB2	27:B1:13:ILE:C	2.31	0.51
32:B6:30:THR:O	32:B6:31:PRO:C	2.49	0.51
34:B8:62:LEU:N	34:B8:63:PRO:HD2	2.26	0.51
35:BA:132:G:C8	35:BA:132:G:H5'	2.41	0.51
35:BA:1382:G:O2'	35:BA:1383:C:H5'	2.10	0.51
35:BA:1424:G:OP1	38:BD:33:LEU:HD21	2.11	0.51
35:BA:1495:A:H2'	35:BA:1495:A:N3	2.26	0.51
35:BA:1771:C:H1'	35:BA:1786:A:C8	2.45	0.51
35:BA:17:G:H2'	35:BA:18:C:C6	2.45	0.51
35:BA:2056:G:N2	35:BA:2057:A:C1'	2.74	0.51
35:BA:2241:A:H2'	35:BA:2242:G:H8	1.73	0.51
35:BA:2064:C:H1'	35:BA:2450:A:C6	2.46	0.51
35:BA:2811:G:H8	35:BA:2811:G:H5''	1.76	0.51
35:BA:2840:C:H2'	35:BA:2841:C:H6	1.76	0.51
35:BA:2886:G:N2	35:BA:2887:U:C2	2.79	0.51
35:BA:290:G:C2'	35:BA:291:C:H5'	2.41	0.51
35:BA:394:A:H2'	35:BA:395:U:H5'	1.91	0.51
35:BA:679:C:O2	35:BA:680:G:C8	2.64	0.51
35:BA:823:G:H2'	35:BA:824:A:C8	2.46	0.51
36:BB:65:C:H2'	36:BB:66:A:H5'	1.92	0.51
38:BD:25:THR:HG23	38:BD:27:THR:HB	1.93	0.51
35:BA:1800:C:OP1	38:BD:264:LYS:NZ	2.44	0.51
38:BD:43:ARG:CB	38:BD:54:ARG:HB2	2.40	0.51
40:BF:117:ARG:CZ	46:BP:5:ASP:N	2.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:80:ALA:HB3	40:BF:83:PHE:CD1	2.46	0.51
42:BH:98:LEU:HD12	42:BH:102:ALA:O	2.11	0.51
42:BH:125:VAL:HG12	42:BH:125:VAL:O	2.11	0.51
42:BH:44:VAL:CG1	42:BH:45:VAL:H	2.23	0.51
45:BO:77:ILE:HD13	50:BT:74:ARG:HG3	1.92	0.51
46:BP:135:LEU:CD1	46:BP:139:LYS:HD2	2.42	0.51
46:BP:21:ARG:HG2	46:BP:21:ARG:O	2.11	0.51
46:BP:48:PRO:CD	46:BP:49:ARG:H	2.24	0.51
47:BQ:59:ARG:HG3	47:BQ:59:ARG:HH11	1.74	0.51
51:BU:55:ARG:HA	51:BU:58:ARG:CG	2.40	0.51
52:BV:61:VAL:CG2	52:BV:99:ILE:HB	2.39	0.51
53:BW:69:LEU:HA	53:BW:109:GLU:HA	1.93	0.51
55:BY:99:CYS:O	55:BY:100:ALA:HB2	2.11	0.51
56:BZ:54:HIS:HB3	56:BZ:101:PRO:HD3	1.93	0.51
1:CA:1419:G:N2	1:CA:1482:G:C1'	2.74	0.51
1:CA:638:G:O2'	1:CA:639:G:H5'	2.11	0.51
1:CA:640:A:O2'	1:CA:641:U:H5'	2.10	0.51
1:CA:819:A:N7	1:CA:1529:G:C6	2.79	0.51
1:CA:882:C:O2'	1:CA:883:C:H5'	2.10	0.51
2:CB:234:PRO:O	2:CB:236:TYR:N	2.44	0.51
1:CA:18:C:C5'	5:CE:127:ASN:HD21	2.24	0.51
5:CE:95:ALA:O	5:CE:96:PRO:C	2.50	0.51
6:CF:30:LEU:HB3	6:CF:35:ALA:HB3	1.92	0.51
7:CG:119:ARG:O	7:CG:120:ILE:C	2.50	0.51
8:CH:104:ARG:O	8:CH:107:LEU:HB3	2.10	0.51
13:CM:82:MET:HB3	13:CM:93:ARG:NH1	2.26	0.51
13:CM:2:ALA:O	13:CM:9:ILE:HG13	2.11	0.51
17:CQ:15:MET:HG2	17:CQ:16:GLN:H	1.75	0.51
18:CR:37:VAL:O	18:CR:38:GLU:C	2.48	0.51
18:CR:66:LEU:HG	18:CR:70:ILE:HD11	1.92	0.51
19:CS:9:VAL:O	19:CS:11:VAL:N	2.44	0.51
19:CS:6:LYS:CD	19:CS:6:LYS:H	2.22	0.51
20:CT:32:ALA:C	20:CT:36:LEU:HD23	2.31	0.51
20:CT:72:LEU:HB3	20:CT:76:ALA:HB1	1.91	0.51
27:D1:16:ASN:N	27:D1:16:ASN:ND2	2.51	0.51
31:D5:11:THR:O	31:D5:12:SER:C	2.49	0.51
32:D6:12:GLU:HB2	32:D6:23:THR:HG22	1.92	0.51
35:DA:1181:C:H2'	35:DA:1182:A:C8	2.46	0.51
35:DA:1465:G:C4'	35:DA:1528:A:H8	2.24	0.51
35:DA:178:G:O2'	35:DA:179:G:H5'	2.11	0.51
35:DA:1863:G:H1	35:DA:1879:C:N4	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1885:A:H3'	35:DA:1886:C:H6	1.74	0.51
35:DA:1952:A:C4	45:DO:22:ILE:HD12	2.46	0.51
35:DA:2303:G:H2'	35:DA:2304:G:O4'	2.11	0.51
35:DA:2313:C:H5'	35:DA:2313:C:C6	2.46	0.51
35:DA:2481:G:HO2'	35:DA:2482:G:P	2.33	0.51
35:DA:2534:A:H2'	35:DA:2535:G:O4'	2.11	0.51
35:DA:558:G:OP1	44:DN:111:PRO:HD2	2.10	0.51
35:DA:585:G:H2'	35:DA:586:A:N7	2.26	0.51
35:DA:631:A:H4'	46:DP:65:ARG:HA	1.93	0.51
36:DB:28:C:O2'	36:DB:29:A:H5'	2.11	0.51
38:DD:54:ARG:H	38:DD:218:ARG:HG3	1.76	0.51
38:DD:21:PHE:O	38:DD:23:GLU:N	2.43	0.51
35:DA:1827:C:OP2	38:DD:222:ARG:NH1	2.44	0.51
38:DD:247:ALA:CA	38:DD:254:THR:HG22	2.39	0.51
40:DF:114:VAL:O	40:DF:117:ARG:N	2.44	0.51
41:DG:104:GLU:C	41:DG:106:LEU:N	2.64	0.51
41:DG:44:GLY:O	41:DG:46:ALA:N	2.44	0.51
43:DI:83:ALA:HA	43:DI:89:TYR:CD1	2.46	0.51
44:DN:27:ALA:O	44:DN:28:THR:C	2.48	0.51
45:DO:105:GLU:HA	45:DO:108:GLU:HG3	1.93	0.51
46:DP:34:GLY:O	46:DP:35:HIS:C	2.49	0.51
47:DQ:9:TYR:CD2	47:DQ:9:TYR:C	2.84	0.51
50:DT:120:ARG:HA	50:DT:123:GLN:HG2	1.92	0.51
50:DT:81:PRO:C	50:DT:82:LEU:HD12	2.31	0.51
35:DA:519:U:H4'	53:DW:25:ARG:HH22	1.75	0.51
1:AA:819:A:N7	1:AA:1529:G:N1	2.58	0.50
1:AA:397:A:H5'	1:AA:398:C:OP1	2.11	0.50
1:AA:512:U:H2'	1:AA:513:C:H6	1.76	0.50
1:AA:599:C:H2'	1:AA:600:C:C6	2.47	0.50
1:AA:665:A:C2'	1:AA:725:G:N2	2.73	0.50
2:AB:81:VAL:HG22	2:AB:215:LEU:HG	1.92	0.50
7:AG:26:PHE:CE1	7:AG:105:VAL:HG22	2.46	0.50
8:AH:85:ARG:HG3	8:AH:85:ARG:HH11	1.76	0.50
9:AI:126:SER:C	9:AI:128:ARG:H	2.14	0.50
9:AI:47:LEU:N	9:AI:47:LEU:HD12	2.24	0.50
11:AK:57:THR:O	11:AK:60:ALA:HB3	2.11	0.50
12:AL:54:LYS:O	12:AL:70:ILE:HD13	2.11	0.50
12:AL:89:ARG:HD3	12:AL:89:ARG:C	2.31	0.50
15:AO:87:ILE:HG22	15:AO:88:ARG:N	2.26	0.50
17:AQ:15:MET:HG2	17:AQ:16:GLN:H	1.76	0.50
18:AR:30:ASP:C	18:AR:32:ARG:H	2.13	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AR:59:SER:OG	18:AR:60:ALA:N	2.44	0.50
13:AM:87:TYR:N	19:AS:73:GLU:O	2.34	0.50
25:AY:78:ALA:HA	25:AY:81:LYS:HG3	1.93	0.50
27:B1:93:GLU:N	27:B1:93:GLU:CD	2.65	0.50
28:B2:51:ARG:O	28:B2:52:ASP:HB2	2.11	0.50
34:B8:5:LYS:HD2	34:B8:5:LYS:N	2.27	0.50
35:BA:1390:U:H6	35:BA:1390:U:OP2	1.93	0.50
35:BA:1669:A:H5''	35:BA:2550:G:OP1	2.10	0.50
35:BA:1885:A:H3'	35:BA:1886:C:C6	2.46	0.50
35:BA:2264:C:H2'	35:BA:2265:U:H6	1.77	0.50
35:BA:2850:A:H5'	35:BA:2868:A:H2	1.76	0.50
35:BA:426:C:C2'	35:BA:427:U:H5'	2.41	0.50
35:BA:604:G:H2'	35:BA:605:C:H6	1.75	0.50
35:BA:619:G:O5'	35:BA:620:G:N2	2.44	0.50
35:BA:662:G:H2'	35:BA:663:G:H8	1.74	0.50
38:BD:118:VAL:CG2	38:BD:119:ALA:N	2.73	0.50
38:BD:24:ILE:HD13	38:BD:24:ILE:O	2.11	0.50
39:BE:34:VAL:O	39:BE:35:GLN:HB2	2.11	0.50
40:BF:3:GLU:HG3	40:BF:19:GLU:HB2	1.92	0.50
41:BG:55:LYS:C	41:BG:57:ALA:H	2.14	0.50
42:BH:130:ARG:NH1	42:BH:130:ARG:HB2	2.27	0.50
43:BI:33:ARG:O	43:BI:35:LEU:HG	2.11	0.50
43:BI:6:LEU:O	43:BI:8:PRO:N	2.44	0.50
45:BO:104:ARG:HG2	45:BO:105:GLU:OE1	2.12	0.50
45:BO:104:ARG:C	45:BO:106:LEU:N	2.59	0.50
46:BP:48:PRO:O	46:BP:49:ARG:C	2.49	0.50
47:BQ:20:ALA:HB2	47:BQ:99:PRO:CD	2.41	0.50
51:BU:38:THR:O	51:BU:41:ALA:HB3	2.12	0.50
35:BA:996:A:O4'	51:BU:92:ARG:NH2	2.44	0.50
52:BV:22:VAL:CG2	52:BV:96:ILE:HB	2.42	0.50
28:B2:22:GLU:HB3	54:BX:5:TYR:CD1	2.46	0.50
54:BX:51:VAL:CG1	54:BX:80:ILE:N	2.75	0.50
55:BY:37:VAL:CG2	55:BY:38:ILE:N	2.71	0.50
55:BY:87:LYS:CG	55:BY:88:LYS:N	2.73	0.50
56:BZ:146:ILE:HG13	56:BZ:147:GLY:N	2.26	0.50
1:CA:1251:A:H2'	1:CA:1252:A:C8	2.46	0.50
1:CA:1409:C:H2'	1:CA:1410:G:C8	2.46	0.50
1:CA:227:G:H2'	1:CA:228:A:C8	2.46	0.50
1:CA:346:G:P	45:DO:107:ARG:HH22	2.33	0.50
1:CA:390:C:H2'	1:CA:391:G:H8	1.76	0.50
1:CA:740:U:O3'	15:CO:39:LEU:HD23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:221:LEU:O	2:CB:221:LEU:HD13	2.10	0.50
4:CD:14:ARG:C	4:CD:16:GLY:H	2.15	0.50
4:CD:153:ARG:CB	4:CD:153:ARG:HH11	2.24	0.50
6:CF:88:VAL:HG12	6:CF:89:MET:N	2.25	0.50
7:CG:145:ALA:O	7:CG:147:ALA:N	2.36	0.50
9:CI:126:SER:C	9:CI:128:ARG:H	2.14	0.50
9:CI:3:GLN:CG	9:CI:20:ARG:HH21	2.24	0.50
11:CK:32:ILE:HD12	11:CK:72:ALA:HB2	1.92	0.50
11:CK:96:ARG:HA	11:CK:99:GLN:CG	2.40	0.50
1:CA:552:U:H4'	12:CL:86:ARG:HG2	1.92	0.50
16:CP:20:VAL:HG22	16:CP:21:VAL:N	2.26	0.50
27:D1:23:LYS:O	27:D1:37:ILE:HG12	2.11	0.50
31:D5:20:ARG:O	31:D5:23:HIS:CD2	2.65	0.50
33:D7:34:ARG:HE	33:D7:39:ARG:HE	1.58	0.50
34:D8:60:LEU:HD23	34:D8:60:LEU:N	2.26	0.50
35:DA:1024:G:N2	35:DA:1142(A):A:H2	2.09	0.50
35:DA:1177:A:H3'	35:DA:1177:A:P	2.50	0.50
35:DA:1357:U:O2'	35:DA:1358:G:H5'	2.11	0.50
35:DA:1487:G:N2	35:DA:1488:G:H1'	2.25	0.50
35:DA:1578:U:C2'	35:DA:1579:A:H5''	2.41	0.50
35:DA:1678:G:C8	35:DA:1678:G:OP2	2.64	0.50
35:DA:576:U:OP1	35:DA:2503:A:OP1	2.30	0.50
35:DA:2758:A:C3'	35:DA:2759:G:H5''	2.41	0.50
35:DA:2777:G:H4'	35:DA:2778:A:H5'	1.92	0.50
35:DA:2830:G:H5'	39:DE:58:ARG:CZ	2.40	0.50
35:DA:305:U:O2'	35:DA:306:U:H5'	2.10	0.50
35:DA:442:G:O4'	40:DF:46:ARG:HG2	2.11	0.50
35:DA:587:C:H3'	46:DP:33:ARG:NH1	2.23	0.50
35:DA:669:G:O2'	35:DA:670:A:OP1	2.27	0.50
35:DA:681:G:H2'	35:DA:682:G:O4'	2.11	0.50
36:DB:56:G:H4'	36:DB:57:A:H8	1.75	0.50
38:DD:112:GLN:N	38:DD:112:GLN:OE1	2.44	0.50
38:DD:12:SER:HB2	38:DD:208:LYS:HB3	1.94	0.50
35:DA:1828:G:O6	38:DD:222:ARG:HD3	2.10	0.50
40:DF:99:TYR:CG	40:DF:99:TYR:O	2.64	0.50
42:DH:85:LYS:O	42:DH:132:ARG:CA	2.58	0.50
44:DN:13:TRP:O	44:DN:135:PRO:HG2	2.11	0.50
44:DN:46:VAL:CG1	44:DN:48:MET:HG3	2.40	0.50
45:DO:119:PRO:HB2	50:DT:68:TYR:HE1	1.72	0.50
46:DP:21:ARG:HG2	46:DP:21:ARG:O	2.09	0.50
47:DQ:132:VAL:HG12	47:DQ:133:ARG:N	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:74:TYR:HB3	47:DQ:91:GLU:CD	2.31	0.50
50:DT:130:ALA:O	50:DT:131:ALA:C	2.49	0.50
50:DT:22:PHE:O	50:DT:23:ARG:HB3	2.11	0.50
51:DU:3:ARG:NH1	51:DU:3:ARG:CG	2.69	0.50
51:DU:66:ASN:OD1	51:DU:76:TYR:HB2	2.11	0.50
51:DU:91:ASP:OD2	51:DU:96:ALA:HB2	2.11	0.50
54:DX:52:VAL:H	54:DX:80:ILE:CG2	2.23	0.50
54:DX:85:PRO:O	54:DX:86:GLY:C	2.47	0.50
56:DZ:89:PHE:HE1	56:DZ:96:VAL:HG21	1.75	0.50
1:AA:1277:C:O2'	1:AA:1278:U:H5'	2.12	0.50
1:AA:1501:C:OP1	1:AA:1508:G:H4'	2.12	0.50
1:AA:1505:G:H4'	1:AA:1506:U:C5'	2.34	0.50
1:AA:277:C:P	17:AQ:41:LYS:HZ1	2.33	0.50
1:AA:540:G:C2	1:AA:541:G:C4	2.99	0.50
1:AA:838:G:O2'	1:AA:839:U:H5''	2.11	0.50
1:AA:957:U:H1'	1:AA:960:U:C4	2.45	0.50
3:AC:171:GLY:O	3:AC:172:ARG:O	2.29	0.50
6:AF:21:LEU:C	6:AF:25:ILE:HG12	2.31	0.50
9:AI:4:TYR:CE1	9:AI:21:PRO:HD3	2.47	0.50
12:AL:89:ARG:HH11	12:AL:89:ARG:CB	2.24	0.50
13:AM:49:THR:HB	13:AM:52:GLU:CG	2.37	0.50
13:AM:69:GLU:HA	13:AM:70:LEU:N	2.25	0.50
15:AO:10:LYS:HG3	15:AO:11:VAL:N	2.26	0.50
15:AO:4:THR:OG1	15:AO:7:GLU:HB2	2.12	0.50
16:AP:58:TYR:O	16:AP:61:SER:OG	2.22	0.50
19:AS:12:ASP:CB	19:AS:15:LEU:HD23	2.41	0.50
20:AT:24:LEU:O	20:AT:27:LYS:HB3	2.11	0.50
7:AG:86:GLN:HG2	23:AW:33:C:C4'	2.41	0.50
25:AY:184:LEU:C	25:AY:184:LEU:HD23	2.31	0.50
27:B1:18:ILE:CG2	27:B1:42:GLN:O	2.58	0.50
28:B2:38:GLN:O	28:B2:39:ALA:C	2.49	0.50
32:B6:20:ASN:ND2	32:B6:21:TYR:N	2.51	0.50
35:BA:1269:A:H2'	35:BA:1270:C:C6	2.46	0.50
35:BA:1707:G:H2'	35:BA:1708:C:H6	1.76	0.50
35:BA:1825:A:H2'	35:BA:1826:G:C8	2.47	0.50
26:B0:41:ARG:HH21	35:BA:2387:U:H1'	1.76	0.50
35:BA:948:G:H2'	35:BA:949:C:H6	1.75	0.50
38:BD:113:VAL:O	38:BD:115:GLN:N	2.32	0.50
38:BD:167:GLY:O	38:BD:168:ARG:HG2	2.12	0.50
38:BD:211:ARG:HG2	38:BD:211:ARG:HH11	1.75	0.50
39:BE:32:PRO:O	39:BE:34:VAL:HG13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:22:ALA:HA	40:BF:26:ALA:HB2	1.92	0.50
42:BH:17:VAL:HB	42:BH:45:VAL:HG22	1.92	0.50
42:BH:89:ILE:CG1	42:BH:90:LYS:N	2.74	0.50
43:BI:109:ILE:O	43:BI:109:ILE:HD12	2.10	0.50
43:BI:72:LEU:O	43:BI:138:ILE:HG12	2.10	0.50
45:BO:80:ASP:HB2	50:BT:71:GLY:O	2.10	0.50
46:BP:106:LEU:C	46:BP:107:LYS:HG2	2.32	0.50
46:BP:39:LYS:C	46:BP:41:ARG:N	2.65	0.50
48:BR:21:TYR:OH	48:BR:43:GLU:HG2	2.11	0.50
48:BR:73:VAL:HG23	48:BR:74:LYS:HD2	1.93	0.50
49:BS:92:TYR:CD2	49:BS:97:ARG:NH1	2.79	0.50
51:BU:68:ALA:O	51:BU:71:GLN:N	2.44	0.50
53:BW:10:VAL:HB	53:BW:101:SER:O	2.11	0.50
56:BZ:137:ILE:HD11	56:BZ:156:LYS:O	2.11	0.50
1:CA:1077:G:N2	1:CA:1079:G:H3'	2.26	0.50
1:CA:979:C:OP1	1:CA:1223:C:N4	2.44	0.50
1:CA:622:A:C8	1:CA:623:C:C6	3.00	0.50
2:CB:118:LEU:HD11	2:CB:141:GLU:OE1	2.10	0.50
2:CB:92:TYR:CD2	2:CB:151:GLY:HA3	2.46	0.50
2:CB:157:ARG:CG	2:CB:158:LEU:H	2.24	0.50
3:CC:172:ARG:HB3	3:CC:174:PRO:HD3	1.94	0.50
3:CC:24:ALA:HB1	3:CC:28:GLN:O	2.10	0.50
3:CC:56:ASP:O	3:CC:57:ILE:HG13	2.10	0.50
5:CE:33:VAL:CG2	5:CE:43:LEU:HD13	2.38	0.50
11:CK:28:THR:CG2	11:CK:29:ILE:N	2.73	0.50
11:CK:48:ILE:N	11:CK:48:ILE:HD13	2.26	0.50
12:CL:20:LYS:H	12:CL:20:LYS:HD3	1.76	0.50
17:CQ:80:GLY:O	17:CQ:81:ARG:HG2	2.11	0.50
17:CQ:86:GLU:HA	17:CQ:89:LEU:HB3	1.93	0.50
19:CS:72:GLY:C	19:CS:74:PHE:H	2.15	0.50
13:CM:91:ARG:HD3	19:CS:81:ARG:NH2	2.26	0.50
23:CW:55:5MU:H73	23:CW:56:U:O4	2.11	0.50
23:CW:9:G:H3'	23:CW:10:G:C8	2.47	0.50
27:D1:68:PRO:O	27:D1:70:VAL:HG23	2.12	0.50
31:D5:30:LEU:HD23	31:D5:41:PRO:CA	2.41	0.50
35:DA:1576:U:H2'	35:DA:1577:C:H6	1.76	0.50
35:DA:1676:A:H2'	35:DA:1677:A:O4'	2.12	0.50
35:DA:1804:C:O2'	35:DA:1805:U:H5'	2.11	0.50
35:DA:184:C:O2'	35:DA:185:U:H5'	2.10	0.50
35:DA:1910:G:C6	35:DA:1921:G:C6	2.98	0.50
35:DA:2545:G:N2	35:DA:2546:U:H1'	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:38:A:H2'	35:DA:39:C:C6	2.46	0.50
36:DB:28:C:N4	36:DB:56:G:H1	2.10	0.50
38:DD:117:VAL:CG2	38:DD:118:VAL:N	2.74	0.50
38:DD:245:PRO:O	38:DD:246:PRO:C	2.48	0.50
39:DE:117:MET:SD	39:DE:136:ARG:NH1	2.85	0.50
39:DE:12:THR:HG23	50:DT:8:LYS:HE2	1.92	0.50
35:DA:2810:A:O2'	39:DE:61:ARG:NE	2.43	0.50
40:DF:184:TYR:HD2	40:DF:185:ASP:N	2.09	0.50
41:DG:120:LEU:H	41:DG:179:PRO:HG2	1.75	0.50
42:DH:85:LYS:HZ1	42:DH:144:VAL:C	2.13	0.50
44:DN:42:TRP:CD2	44:DN:44:PRO:HD3	2.45	0.50
46:DP:96:THR:HB	46:DP:97:PRO:CD	2.41	0.50
47:DQ:88:GLY:C	47:DQ:90:VAL:H	2.12	0.50
48:DR:116:LEU:O	48:DR:117:VAL:CB	2.58	0.50
50:DT:23:ARG:HB2	50:DT:24:PRO:HD2	1.93	0.50
55:DY:98:VAL:O	55:DY:99:CYS:SG	2.68	0.50
56:DZ:137:ILE:CD1	56:DZ:157:LEU:HD23	2.41	0.50
1:AA:15:G:H4'	5:AE:24:ARG:HH21	1.77	0.50
1:AA:167:G:O2'	1:AA:168:G:H5'	2.11	0.50
1:AA:189(B):C:H2'	1:AA:189(C):C:C6	2.46	0.50
1:AA:255:G:H2'	1:AA:256:U:C6	2.45	0.50
1:AA:740:U:O3'	15:AO:39:LEU:HD23	2.11	0.50
5:AE:121:LYS:HD2	5:AE:122:GLU:H	1.77	0.50
6:AF:43:LEU:N	6:AF:43:LEU:HD12	2.26	0.50
7:AG:103:TRP:O	7:AG:104:LEU:C	2.49	0.50
7:AG:50:ILE:HG13	7:AG:58:PRO:HB3	1.94	0.50
1:AA:1368:G:OP2	9:AI:112:LYS:HE3	2.11	0.50
10:AJ:96:ILE:HD13	10:AJ:96:ILE:N	2.26	0.50
11:AK:69:ALA:O	11:AK:70:LYS:C	2.50	0.50
13:AM:74:VAL:O	13:AM:78:ILE:HG13	2.11	0.50
14:AN:36:PHE:CD1	14:AN:36:PHE:O	2.64	0.50
15:AO:64:ARG:O	15:AO:65:ARG:C	2.49	0.50
16:AP:28:ARG:HG2	16:AP:28:ARG:NH1	2.24	0.50
16:AP:6:LEU:HD23	16:AP:17:TYR:CD2	2.47	0.50
16:AP:82:GLN:O	16:AP:84:ALA:N	2.45	0.50
17:AQ:40:LYS:HG2	17:AQ:41:LYS:N	2.26	0.50
18:AR:26:LEU:HD21	18:AR:42:ARG:CZ	2.41	0.50
19:AS:20:LEU:O	19:AS:23:ASN:HB3	2.11	0.50
25:AY:63:PRO:HB2	25:AY:64:ARG:NH2	2.27	0.50
28:B2:55:ARG:C	28:B2:57:ILE:H	2.14	0.50
35:BA:1195:G:H2'	35:BA:1196:C:C6	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1714:G:H2'	35:BA:1717:G:H8	1.75	0.50
35:BA:2025:C:H2'	35:BA:2026:C:C6	2.46	0.50
35:BA:2065:C:H2'	35:BA:2066:C:H6	1.76	0.50
35:BA:2221:G:H5'	35:BA:2222:G:OP2	2.12	0.50
35:BA:2597:G:H2'	35:BA:2598:A:C8	2.47	0.50
35:BA:2759:G:H5'	35:BA:2759:G:C8	2.46	0.50
35:BA:382:G:H1	35:BA:392:C:H42	1.58	0.50
35:BA:447:A:C4	35:BA:473:G:N7	2.80	0.50
35:BA:501:A:H2'	35:BA:502:A:C8	2.46	0.50
35:BA:261:G:C1'	35:BA:609:A:H2	2.24	0.50
35:BA:614:U:H4'	35:BA:614(C):A:H62	1.76	0.50
35:BA:779:U:H2'	35:BA:780:G:H8	1.76	0.50
37:BC:89:ALA:HA	37:BC:153:ILE:CB	2.40	0.50
38:BD:176:ARG:HG2	38:BD:176:ARG:NH1	2.26	0.50
39:BE:132:HIS:O	39:BE:135:HIS:CD2	2.64	0.50
39:BE:173:VAL:O	39:BE:174:ASP:C	2.49	0.50
40:BF:7:TYR:HB3	40:BF:16:GLY:O	2.11	0.50
43:BI:2:LYS:O	43:BI:3:VAL:HG13	2.11	0.50
44:BN:126:PRO:O	44:BN:127:ASP:CB	2.60	0.50
45:BO:68:GLU:N	45:BO:68:GLU:OE2	2.44	0.50
46:BP:68:GLN:HG3	46:BP:68:GLN:O	2.11	0.50
35:BA:1278:A:C5'	48:BR:36:THR:HG22	2.42	0.50
50:BT:100:TYR:CD1	50:BT:100:TYR:N	2.79	0.50
50:BT:108:ARG:NH1	50:BT:108:ARG:HB3	2.27	0.50
45:BO:104:ARG:NH2	50:BT:33:LYS:HE3	2.26	0.50
52:BV:1:MET:CE	52:BV:45:THR:H	2.24	0.50
54:BX:31:HIS:ND1	54:BX:32:PRO:HD2	2.26	0.50
55:BY:2:ARG:N	55:BY:5:MET:CE	2.74	0.50
56:BZ:58:VAL:HA	56:BZ:68:PRO:CA	2.41	0.50
56:BZ:3:TYR:O	56:BZ:58:VAL:N	2.44	0.50
47:BQ:141:GLN:HG2	56:BZ:72:ARG:HG2	1.93	0.50
1:CA:1486:G:H2'	1:CA:1487:G:C8	2.47	0.50
1:CA:940:C:H2'	1:CA:941:G:C8	2.47	0.50
1:CA:989:C:O2'	1:CA:990:C:H5'	2.11	0.50
3:CC:127:ARG:NH1	3:CC:127:ARG:HG2	2.23	0.50
3:CC:84:ILE:CA	3:CC:87:LEU:HD12	2.38	0.50
4:CD:128:VAL:CG1	4:CD:129:ASN:H	2.14	0.50
6:CF:85:VAL:HG12	6:CF:85:VAL:O	2.11	0.50
7:CG:122:HIS:O	7:CG:125:MET:N	2.44	0.50
7:CG:152:ALA:C	7:CG:154:TYR:N	2.65	0.50
9:CI:102:LEU:O	9:CI:103:THR:OG1	2.22	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:76:ASN:O	12:CL:77:LEU:HD23	2.10	0.50
12:CL:89:ARG:NE	12:CL:91:LYS:HE2	2.27	0.50
15:CO:4:THR:OG1	15:CO:7:GLU:HB2	2.12	0.50
25:CY:64:ARG:HA	25:CY:103:ILE:HG13	1.91	0.50
35:DA:818:G:N2	35:DA:1190:G:C6	2.80	0.50
35:DA:1215:G:H2'	35:DA:1216:G:O4'	2.11	0.50
35:DA:1348:G:H2'	35:DA:1349:A:C5'	2.28	0.50
35:DA:1495:A:N3	35:DA:1495:A:H2'	2.26	0.50
35:DA:1649:G:O2'	35:DA:1650:G:H5'	2.11	0.50
35:DA:1682:G:O2'	35:DA:1683:C:H5'	2.11	0.50
35:DA:1991:U:C2'	35:DA:1992:G:C5'	2.90	0.50
35:DA:2368:C:H2'	35:DA:2369:A:C8	2.43	0.50
35:DA:2639:A:C2'	35:DA:2640:G:C5'	2.82	0.50
35:DA:566:U:H2'	35:DA:567:A:C8	2.46	0.50
35:DA:939:G:O2'	35:DA:940:G:H5'	2.11	0.50
35:DA:956:G:H22	35:DA:959:A:H3'	1.75	0.50
38:DD:111:LEU:HD13	38:DD:112:GLN:N	2.26	0.50
38:DD:111:LEU:HD22	38:DD:115:GLN:OE1	2.12	0.50
38:DD:223:GLY:HA2	38:DD:226:MET:SD	2.52	0.50
35:DA:2591:C:P	38:DD:239:ARG:HG3	2.51	0.50
38:DD:35:LYS:CG	38:DD:64:ILE:H	2.10	0.50
39:DE:167:VAL:CG2	39:DE:168:MET:H	2.21	0.50
39:DE:197:ILE:CG1	39:DE:199:ARG:HH12	2.22	0.50
40:DF:185:ASP:HA	40:DF:188:ARG:CB	2.40	0.50
40:DF:125:LEU:HD12	40:DF:196:LEU:CD2	2.41	0.50
41:DG:114:ILE:CG2	41:DG:115:ARG:H	2.17	0.50
41:DG:46:ALA:HA	41:DG:51:ARG:HD3	1.93	0.50
42:DH:140:LYS:O	42:DH:144:VAL:HG23	2.11	0.50
43:DI:102:SER:HB2	43:DI:109:ILE:CG2	2.40	0.50
43:DI:45:LYS:O	43:DI:48:GLU:HB3	2.12	0.50
46:DP:34:GLY:O	46:DP:36:LYS:HG3	2.11	0.50
46:DP:48:PRO:O	46:DP:49:ARG:C	2.49	0.50
47:DQ:57:HIS:C	47:DQ:57:HIS:ND1	2.64	0.50
48:DR:35:THR:HG23	48:DR:112:ALA:O	2.11	0.50
49:DS:66:ALA:HA	49:DS:69:VAL:CG1	2.41	0.50
49:DS:90:GLY:HA2	49:DS:92:TYR:CE2	2.47	0.50
51:DU:88:ILE:O	51:DU:90:VAL:N	2.44	0.50
51:DU:98:LEU:C	51:DU:100:VAL:H	2.14	0.50
52:DV:18:LEU:CD2	52:DV:19:LYS:H	2.11	0.50
52:DV:1:MET:HE1	52:DV:46:VAL:HG23	1.93	0.50
52:DV:88:ARG:HG3	52:DV:88:ARG:NH1	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DV:98:GLU:N	52:DV:98:GLU:OE1	2.44	0.50
35:DA:456:C:N3	54:DX:66:LEU:HD22	2.26	0.50
56:DZ:47:VAL:HG12	56:DZ:51:ALA:HB2	1.93	0.50
1:AA:1029:C:H1'	1:AA:1033:G:N1	2.26	0.50
1:AA:1251:A:O2'	1:AA:1252:A:H5'	2.12	0.50
1:AA:199:G:H2'	1:AA:200:G:C8	2.45	0.50
1:AA:251:G:N2	1:AA:253:U:C4	2.80	0.50
1:AA:429:U:C1'	1:AA:430:A:H5''	2.41	0.50
1:AA:658:G:O2'	1:AA:659:U:H5'	2.10	0.50
1:AA:90:U:H6	1:AA:90:U:O5'	1.94	0.50
1:AA:973:G:O4'	10:AJ:55:LYS:HG2	2.11	0.50
2:AB:196:LEU:O	2:AB:196:LEU:HD12	2.11	0.50
2:AB:58:ILE:O	2:AB:61:LEU:HB3	2.11	0.50
4:AD:101:LEU:HD23	4:AD:121:VAL:CG1	2.40	0.50
4:AD:119:GLN:NE2	4:AD:123:HIS:NE2	2.59	0.50
4:AD:29:PRO:C	4:AD:30:LYS:HG2	2.32	0.50
4:AD:96:LEU:HD13	4:AD:96:LEU:N	2.27	0.50
7:AG:152:ALA:C	7:AG:154:TYR:N	2.65	0.50
7:AG:50:ILE:HG21	7:AG:61:VAL:HG21	1.93	0.50
7:AG:66:VAL:O	7:AG:69:VAL:N	2.39	0.50
8:AH:4:ASP:OD2	8:AH:7:ALA:N	2.33	0.50
8:AH:63:LEU:HG	8:AH:65:TYR:OH	2.11	0.50
9:AI:4:TYR:HD1	9:AI:4:TYR:N	2.09	0.50
12:AL:46:LYS:CG	12:AL:47:LYS:N	2.70	0.50
15:AO:9:GLN:O	15:AO:10:LYS:C	2.50	0.50
23:AW:32:G:N3	23:AW:32:G:H2'	2.26	0.50
23:AW:8:U:O2'	23:AW:47:G:N2	2.45	0.50
25:AY:118:VAL:C	25:AY:120:GLN:N	2.64	0.50
29:B3:23:LEU:O	29:B3:28:LEU:HB2	2.10	0.50
29:B3:28:LEU:HA	29:B3:33:GLN:OE1	2.12	0.50
30:B4:43:TYR:C	30:B4:45:GLY:H	2.14	0.50
35:BA:1267:U:H2'	35:BA:1267:U:O2	2.11	0.50
35:BA:1411:C:O2'	35:BA:1412:A:H8	1.93	0.50
35:BA:1563:G:O2'	35:BA:1564:C:H5'	2.12	0.50
35:BA:1848:A:H2'	35:BA:1849:G:H8	1.75	0.50
35:BA:1964:G:H3'	35:BA:1965:C:C5'	2.42	0.50
35:BA:1986:A:C2'	35:BA:1987:G:H5''	2.41	0.50
35:BA:2119:A:C3'	35:BA:2120:G:H5''	2.42	0.50
35:BA:2425:A:O4'	35:BA:2427:C:C6	2.65	0.50
35:BA:2805:G:H22	35:BA:2893:G:H1	1.59	0.50
35:BA:2821:A:P	48:BR:2:ARG:NH2	2.62	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:992:C:O3'	52:BV:75:PHE:CE2	2.64	0.50
38:BD:130:ALA:CB	38:BD:192:THR:HA	2.42	0.50
38:BD:267:SER:O	38:BD:269:PHE:CD1	2.65	0.50
42:BH:103:LEU:HG	42:BH:105:LEU:CD1	2.42	0.50
42:BH:103:LEU:HD22	42:BH:123:PHE:CE2	2.46	0.50
42:BH:85:LYS:HB3	42:BH:133:VAL:O	2.12	0.50
42:BH:149:ARG:CB	42:BH:162:ILE:HD11	2.41	0.50
42:BH:30:LYS:NZ	42:BH:81:GLU:HA	2.27	0.50
44:BN:42:TRP:CH2	44:BN:44:PRO:HA	2.46	0.50
46:BP:124:LYS:HA	46:BP:143:GLY:HA3	1.92	0.50
47:BQ:28:ALA:HB3	47:BQ:105:GLU:OE2	2.11	0.50
49:BS:79:ALA:C	49:BS:80:LEU:HD12	2.31	0.50
49:BS:90:GLY:HA2	49:BS:92:TYR:CE2	2.46	0.50
50:BT:114:LEU:N	50:BT:114:LEU:HD23	2.26	0.50
50:BT:130:ALA:O	50:BT:131:ALA:C	2.49	0.50
50:BT:53:ARG:NE	50:BT:60:THR:OG1	2.44	0.50
52:BV:4:ILE:HB	52:BV:40:LEU:HD11	1.93	0.50
52:BV:5:VAL:HB	52:BV:37:VAL:O	2.11	0.50
53:BW:66:GLU:HA	53:BW:69:LEU:HD21	1.94	0.50
55:BY:16:ALA:CA	55:BY:21:LYS:HD2	2.42	0.50
55:BY:28:LYS:HA	55:BY:38:ILE:HG22	1.93	0.50
56:BZ:67:LEU:O	56:BZ:69:THR:N	2.43	0.50
1:CA:1226:C:OP1	19:CS:81:ARG:NH1	2.45	0.50
1:CA:1483:A:H3'	1:CA:1484:C:H6	1.75	0.50
1:CA:678:U:H2'	1:CA:679:C:H6	1.72	0.50
1:CA:932:C:H5"	7:CG:3:ARG:CD	2.35	0.50
2:CB:21:ARG:O	2:CB:23:ARG:N	2.44	0.50
2:CB:9:GLU:O	2:CB:13:ALA:CB	2.60	0.50
3:CC:182:ILE:HG23	3:CC:202:ILE:O	2.11	0.50
6:CF:72:VAL:CG1	6:CF:73:ASN:H	2.21	0.50
6:CF:75:LEU:HD23	6:CF:79:LEU:HD21	1.93	0.50
7:CG:27:ILE:HD11	7:CG:43:PHE:CD2	2.46	0.50
1:CA:972:C:C2'	10:CJ:55:LYS:HD3	2.42	0.50
12:CL:9:GLN:O	12:CL:10:LEU:C	2.48	0.50
15:CO:67:LEU:HD22	15:CO:78:TYR:HE1	1.76	0.50
1:CA:279:A:OP2	17:CQ:95:TYR:OH	2.30	0.50
25:CY:171:LYS:O	25:CY:175:LEU:HB2	2.11	0.50
35:DA:1006:C:N3	35:DA:1138:G:C2	2.79	0.50
35:DA:990:A:C6	35:DA:1186:G:H1'	2.46	0.50
35:DA:1433:U:H3	35:DA:1560:G:H1	1.59	0.50
35:DA:1613:G:C6	35:DA:1619:G:C6	2.99	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1778:U:H5	35:DA:1784:A:N3	2.09	0.50
35:DA:1990:C:H2'	35:DA:1991:U:O4'	2.12	0.50
35:DA:2247:A:H2'	35:DA:2248:C:H6	1.76	0.50
35:DA:2280:G:N3	35:DA:2388:A:H2	2.08	0.50
35:DA:2298:A:H62	35:DA:2318:G:H8	1.57	0.50
35:DA:426:C:C2'	35:DA:427:U:H5'	2.41	0.50
35:DA:435:C:H2'	35:DA:436:C:H5'	1.94	0.50
35:DA:778:G:C5	35:DA:779:U:C4	3.00	0.50
35:DA:948:G:H2'	35:DA:949:C:H6	1.77	0.50
37:DC:203:GLY:O	37:DC:204:ALA:HB2	2.12	0.50
37:DC:74:VAL:H	37:DC:91:ALA:HB1	1.77	0.50
35:DA:1792:G:OP2	38:DD:206:LEU:HD12	2.11	0.50
38:DD:260:ARG:HG2	38:DD:260:ARG:NH1	2.26	0.50
38:DD:27:THR:O	38:DD:29:PRO:HD2	2.11	0.50
39:DE:8:LYS:HE3	39:DE:188:VAL:HG13	1.93	0.50
39:DE:9:VAL:CG1	39:DE:25:VAL:HG12	2.40	0.50
39:DE:52:LEU:HB3	39:DE:76:ARG:H	1.76	0.50
40:DF:132:VAL:HG13	40:DF:133:ASN:N	2.27	0.50
43:DI:94:ALA:HA	43:DI:97:ILE:CB	2.42	0.50
44:DN:70:LYS:HG3	44:DN:72:TYR:HE1	1.76	0.50
35:DA:2726:U:H6	45:DO:67:LYS:HZ3	1.58	0.50
45:DO:86:ILE:O	45:DO:93:PRO:HA	2.11	0.50
46:DP:138:LEU:CD2	46:DP:142:GLY:HA3	2.41	0.50
46:DP:70:GLN:CG	46:DP:71:VAL:H	2.24	0.50
48:DR:16:HIS:O	48:DR:19:ALA:N	2.45	0.50
48:DR:82:GLU:OE1	48:DR:83:ILE:HD13	2.11	0.50
49:DS:92:TYR:HD2	49:DS:97:ARG:CZ	2.25	0.50
51:DU:40:PHE:HD2	51:DU:40:PHE:N	2.10	0.50
51:DU:98:LEU:O	51:DU:100:VAL:N	2.44	0.50
52:DV:85:LYS:O	52:DV:87:HIS:N	2.41	0.50
54:DX:73:ARG:O	54:DX:73:ARG:HG3	2.11	0.50
1:AA:112:G:C2'	1:AA:113:G:H5'	2.41	0.50
1:AA:1132:C:O2'	1:AA:1133:G:H5'	2.11	0.50
1:AA:119:A:H4'	1:AA:120:A:O5'	2.12	0.50
1:AA:1222:G:H5''	19:AS:78:ARG:HH11	1.74	0.50
1:AA:1497:G:C2'	1:AA:1498:U:H5'	2.41	0.50
1:AA:1517:G:H1'	35:BA:1919:A:O3'	2.12	0.50
1:AA:183:G:H1	1:AA:194:C:H42	1.59	0.50
1:AA:25:C:O2'	1:AA:26:A:H5'	2.11	0.50
1:AA:358:U:H2'	1:AA:359:U:H6	1.76	0.50
1:AA:714:G:H2'	1:AA:715:A:C8	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:865:A:H2'	1:AA:866:C:H6	1.74	0.50
1:AA:963:G:H2'	1:AA:964:A:H8	1.75	0.50
2:AB:135:GLN:O	2:AB:139:LYS:HB2	2.12	0.50
2:AB:17:PHE:C	2:AB:17:PHE:CD2	2.83	0.50
3:AC:19:GLU:O	3:AC:56:ASP:HA	2.11	0.50
4:AD:100:ARG:O	4:AD:103:ASN:HB3	2.12	0.50
6:AF:28:ARG:NH1	6:AF:28:ARG:HG3	2.27	0.50
9:AI:28:VAL:HA	9:AI:63:ILE:O	2.11	0.50
11:AK:30:VAL:CG2	11:AK:68:ALA:HB2	2.42	0.50
14:AN:4:LYS:HA	14:AN:7:ILE:CD1	2.42	0.50
17:AQ:48:GLU:O	17:AQ:50:LYS:N	2.44	0.50
19:AS:18:LYS:HA	19:AS:21:GLU:HG2	1.92	0.50
23:AW:27:G:H2'	23:AW:28:U:C6	2.46	0.50
26:B0:32:ARG:N	26:B0:35:ASN:HD22	1.89	0.50
26:B0:43:THR:O	26:B0:43:THR:HG23	2.12	0.50
35:BA:1430:C:H2'	35:BA:1431:U:H6	1.77	0.50
35:BA:1786:A:C6	35:BA:1938:A:N7	2.79	0.50
35:BA:1860:G:H2'	35:BA:1861:G:H8	1.77	0.50
35:BA:1956:U:H2'	35:BA:1957:C:H5'	1.93	0.50
35:BA:215:G:H4'	35:BA:216:A:O5'	2.11	0.50
35:BA:2321:G:H2'	35:BA:2321:G:N3	2.26	0.50
35:BA:2441:C:H2'	35:BA:2441:C:O2	2.10	0.50
35:BA:2559:C:H2'	35:BA:2559:C:O2	2.12	0.50
35:BA:2545:G:N3	35:BA:2565:A:H2	2.09	0.50
35:BA:307:G:H21	35:BA:330:A:H62	1.59	0.50
35:BA:366:C:H5''	35:BA:403:U:H3	1.76	0.50
35:BA:703:U:H2'	35:BA:704:G:C5'	2.41	0.50
37:BC:64:LEU:HD12	37:BC:66:HIS:HB2	1.94	0.50
38:BD:239:ARG:NH2	38:BD:239:ARG:HG3	2.21	0.50
38:BD:94:LEU:HD12	38:BD:94:LEU:N	2.14	0.50
39:BE:113:PHE:CE2	39:BE:158:GLY:HA2	2.47	0.50
39:BE:116:VAL:HG22	39:BE:117:MET:H	1.74	0.50
42:BH:92:ILE:HG22	42:BH:93:GLY:H	1.76	0.50
43:BI:88:ILE:HG13	43:BI:122:GLU:HA	1.94	0.50
46:BP:128:HIS:O	46:BP:129:ALA:HB2	2.11	0.50
47:BQ:57:HIS:C	47:BQ:57:HIS:ND1	2.64	0.50
35:BA:869:G:H1'	47:BQ:8:LYS:HZ3	1.74	0.50
49:BS:17:ARG:NE	49:BS:89:ARG:NH2	2.60	0.50
49:BS:17:ARG:HG3	49:BS:18:ILE:HD13	1.92	0.50
49:BS:26:LEU:HD23	49:BS:28:VAL:HG22	1.92	0.50
49:BS:34:HIS:CE1	49:BS:55:ALA:HB2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BS:92:TYR:HD2	49:BS:97:ARG:CZ	2.24	0.50
51:BU:76:TYR:CZ	51:BU:80:ILE:HG13	2.46	0.50
52:BV:52:VAL:C	52:BV:54:GLY:N	2.64	0.50
55:BY:96:ILE:HG13	55:BY:100:ALA:H	1.75	0.50
55:BY:101:LYS:CG	55:BY:102:CYS:N	2.72	0.50
47:BQ:141:GLN:HE21	56:BZ:72:ARG:HG2	1.76	0.50
1:CA:1287:A:H2	1:CA:1353:G:N3	2.10	0.50
1:CA:1333:A:H3'	1:CA:1334:G:H8	1.76	0.50
1:CA:380:G:N1	1:CA:384:G:C6	2.79	0.50
1:CA:441:A:C6	1:CA:494:U:C2	2.99	0.50
1:CA:473:G:H2'	1:CA:474:G:C8	2.45	0.50
1:CA:300:A:H2	1:CA:566:G:O6	1.94	0.50
1:CA:658:G:H1'	15:CO:22:THR:CB	2.38	0.50
1:CA:740:U:O2'	1:CA:741:G:H5'	2.10	0.50
2:CB:196:LEU:HD12	2:CB:196:LEU:O	2.11	0.50
4:CD:173:TRP:O	4:CD:186:LEU:HB2	2.12	0.50
4:CD:56:VAL:C	4:CD:58:LEU:N	2.64	0.50
5:CE:41:VAL:HG13	5:CE:113:ALA:HA	1.94	0.50
5:CE:47:LYS:HD3	5:CE:47:LYS:N	2.26	0.50
5:CE:80:ILE:HD11	5:CE:91:LEU:HB2	1.94	0.50
9:CI:82:ALA:O	9:CI:96:LEU:HD21	2.12	0.50
9:CI:97:LYS:HB3	9:CI:98:PRO:HD3	1.94	0.50
20:CT:29:LYS:O	20:CT:32:ALA:HB3	2.10	0.50
28:D2:56:GLN:HE21	28:D2:56:GLN:HA	1.72	0.50
30:D4:17:GLY:O	30:D4:19:GLY:N	2.41	0.50
31:D5:7:PRO:HA	35:DA:2615:U:C2	2.47	0.50
34:D8:29:LYS:HZ3	34:D8:44:LYS:HB2	1.76	0.50
35:DA:1233:C:C2'	35:DA:1234:U:H5'	2.42	0.50
35:DA:1299:G:H5''	35:DA:1300:U:OP1	2.11	0.50
33:D7:9:ARG:NH1	35:DA:1310:G:OP2	2.44	0.50
35:DA:1720:U:H2'	35:DA:1721:G:H5'	1.92	0.50
35:DA:1819:A:O2'	35:DA:1820:U:OP2	2.24	0.50
1:CA:1409:C:H4'	35:DA:1915:U:O4	2.11	0.50
35:DA:2050:C:H2'	35:DA:2051:A:O4'	2.12	0.50
35:DA:271(H):G:O2'	35:DA:271(I):G:H8	1.94	0.50
35:DA:272(B):G:H2'	35:DA:272(C):G:H8	1.77	0.50
35:DA:694:U:C2'	35:DA:695:G:O5'	2.59	0.50
35:DA:694:U:H2'	35:DA:695:G:O5'	2.11	0.50
38:DD:186:HIS:CD2	38:DD:188:GLU:HG2	2.46	0.50
35:DA:2574:G:O2'	39:DE:143:ASN:HB3	2.12	0.50
39:DE:199:ARG:NH1	39:DE:199:ARG:HG3	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DE:31:CYS:HB3	39:DE:49:LEU:HB3	1.94	0.50
40:DF:28:ILE:HA	40:DF:112:MET:HG2	1.94	0.50
40:DF:32:LEU:C	40:DF:32:LEU:CD2	2.80	0.50
35:DA:2745:C:H1'	42:DH:143:GLN:HG2	1.93	0.50
44:DN:54:VAL:HB	44:DN:122:VAL:HG22	1.94	0.50
44:DN:72:TYR:N	44:DN:85:ILE:O	2.41	0.50
44:DN:87:LEU:O	44:DN:88:GLU:C	2.50	0.50
46:DP:48:PRO:CD	46:DP:49:ARG:H	2.24	0.50
46:DP:59:LEU:CA	46:DP:61:ARG:CZ	2.88	0.50
47:DQ:50:ALA:O	47:DQ:54:MET:HB2	2.11	0.50
50:DT:13:ARG:NH1	50:DT:15:VAL:HG12	2.26	0.50
45:DO:104:ARG:NH2	50:DT:33:LYS:HE3	2.27	0.50
51:DU:7:GLY:O	51:DU:8:VAL:HG22	2.11	0.50
52:DV:2:PHE:HB2	52:DV:42:GLY:O	2.12	0.50
52:DV:66:ARG:HB2	52:DV:95:LEU:H	1.76	0.50
54:DX:12:VAL:HG13	54:DX:17:ALA:HB1	1.94	0.50
54:DX:57:LEU:O	54:DX:58:HIS:CG	2.64	0.50
55:DY:29:GLU:N	55:DY:29:GLU:OE1	2.43	0.50
55:DY:37:VAL:CG2	55:DY:38:ILE:H	2.04	0.50
56:DZ:141:VAL:HG22	56:DZ:141:VAL:O	2.12	0.50
56:DZ:169:GLU:HG2	56:DZ:170:THR:H	1.76	0.50
56:DZ:115:GLY:N	56:DZ:177:PRO:HD3	2.26	0.50
1:AA:1090:U:O2'	1:AA:1091:U:H5'	2.10	0.50
1:AA:1379:G:C4	1:AA:1380:U:H5	2.30	0.50
1:AA:324:G:N2	1:AA:326:G:H3'	2.27	0.50
1:AA:677:U:H2'	1:AA:678:U:C6	2.46	0.50
1:AA:728:A:C2	1:AA:729:A:C5	2.99	0.50
2:AB:101:MET:O	2:AB:105:PHE:CA	2.57	0.50
3:AC:182:ILE:HG23	3:AC:202:ILE:O	2.12	0.50
3:AC:52:LEU:CD2	3:AC:52:LEU:H	2.20	0.50
4:AD:128:VAL:O	4:AD:130:GLY:N	2.41	0.50
4:AD:170:VAL:HG13	4:AD:174:LEU:HB2	1.94	0.50
5:AE:41:VAL:O	5:AE:67:VAL:N	2.44	0.50
5:AE:47:LYS:HD3	5:AE:47:LYS:N	2.26	0.50
7:AG:29:LYS:HG3	7:AG:101:LEU:HB3	1.93	0.50
7:AG:143:ARG:NH1	7:AG:143:ARG:HB2	2.27	0.50
8:AH:107:LEU:C	8:AH:107:LEU:HD23	2.31	0.50
9:AI:118:LYS:H	9:AI:121:ARG:HB3	1.76	0.50
11:AK:32:ILE:HD12	11:AK:72:ALA:HB2	1.94	0.50
11:AK:96:ARG:HA	11:AK:99:GLN:CG	2.36	0.50
13:AM:108:ARG:NH2	13:AM:114:ARG:HG2	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:112:GLY:C	13:AM:113:PRO:HG2	2.32	0.50
13:AM:13:LYS:HB2	13:AM:18:ALA:HB2	1.93	0.50
15:AO:85:LEU:HB2	15:AO:87:ILE:HD11	1.93	0.50
18:AR:25:THR:O	18:AR:25:THR:HG22	2.11	0.50
32:B6:12:GLU:HB2	32:B6:23:THR:HG22	1.94	0.50
35:BA:1177:A:H3'	35:BA:1177:A:P	2.51	0.50
35:BA:1543:C:O2	35:BA:1543:C:C2'	2.60	0.50
35:BA:1652:A:N6	35:BA:1653:G:N2	2.60	0.50
35:BA:1948:G:H2'	35:BA:1949:G:H8	1.76	0.50
35:BA:224:G:O2'	35:BA:225:A:H5'	2.11	0.50
35:BA:2639:A:H3'	35:BA:2640:G:H5'	1.93	0.50
35:BA:2697:G:C2	35:BA:2711:A:C2	2.99	0.50
35:BA:2699:C:O2'	35:BA:2700:C:H5'	2.10	0.50
35:BA:289:A:H3'	35:BA:290:G:H8	1.75	0.50
35:BA:577:G:H2'	35:BA:578:A:H8	1.76	0.50
35:BA:587:C:C5	35:BA:671:C:H1'	2.47	0.50
36:BB:29:A:H2'	36:BB:30:C:C6	2.47	0.50
38:BD:40:THR:HG22	38:BD:41:GLY:O	2.12	0.50
39:BE:152:LYS:HZ3	44:BN:78:TYR:CB	2.24	0.50
39:BE:16:ARG:O	39:BE:17:ASP:HB3	2.11	0.50
39:BE:201:THR:CG2	39:BE:202:LYS:N	2.75	0.50
39:BE:60:ASN:O	39:BE:61:ARG:C	2.50	0.50
40:BF:168:ARG:HA	40:BF:175:THR:HG21	1.92	0.50
40:BF:32:LEU:O	40:BF:33:LEU:C	2.50	0.50
42:BH:85:LYS:HZ1	42:BH:144:VAL:C	2.15	0.50
44:BN:13:TRP:O	44:BN:135:PRO:HG2	2.12	0.50
45:BO:12:ASP:N	45:BO:12:ASP:OD2	2.43	0.50
45:BO:1:MET:CG	45:BO:32:TYR:HD2	2.23	0.50
46:BP:19:VAL:HG23	46:BP:19:VAL:O	2.11	0.50
47:BQ:32:TYR:CD1	47:BQ:32:TYR:N	2.80	0.50
47:BQ:73:PRO:HG3	47:BQ:93:TYR:CD2	2.47	0.50
48:BR:10:LEU:HB3	48:BR:17:ARG:NE	2.27	0.50
48:BR:26:LYS:NZ	48:BR:71:GLN:HB3	2.27	0.50
45:BO:104:ARG:CZ	50:BT:33:LYS:HD2	2.42	0.50
31:B5:20:ARG:NH1	53:BW:15:ARG:NH2	2.53	0.50
53:BW:20:VAL:O	53:BW:23:LEU:N	2.45	0.50
54:BX:25:LYS:O	54:BX:26:TYR:O	2.29	0.50
54:BX:37:THR:C	54:BX:39:ILE:H	2.14	0.50
56:BZ:166:SER:CB	56:BZ:167:PRO:CA	2.86	0.50
56:BZ:42:VAL:HG13	56:BZ:43:GLU:N	2.26	0.50
1:CA:1069:C:O2'	1:CA:1192:C:H1'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1416:G:N1	1:CA:1485:U:O2	2.44	0.50
1:CA:285:G:O2'	1:CA:286:G:H5'	2.11	0.50
1:CA:439:A:C4	1:CA:496:A:C2	3.00	0.50
1:CA:706:A:C5	1:CA:707:C:C5	2.94	0.50
1:CA:854:G:OP2	1:CA:871:U:C5	2.65	0.50
1:CA:939:G:H2'	1:CA:940:C:H6	1.76	0.50
1:CA:963:G:H2'	1:CA:964:A:H8	1.75	0.50
2:CB:80:ILE:HG12	2:CB:211:ILE:HG21	1.93	0.50
2:CB:74:LYS:HG3	2:CB:77:ALA:HB3	1.93	0.50
3:CC:180:ALA:O	3:CC:205:GLY:O	2.29	0.50
3:CC:69:HIS:CD2	3:CC:69:HIS:N	2.78	0.50
8:CH:50:ARG:HH11	8:CH:50:ARG:HG2	1.77	0.50
8:CH:4:ASP:OD1	8:CH:6:ILE:HB	2.11	0.50
1:CA:1250:A:H4'	9:CI:68:GLY:O	2.12	0.50
13:CM:49:THR:HB	13:CM:52:GLU:CG	2.37	0.50
15:CO:67:LEU:HD22	15:CO:78:TYR:CE1	2.47	0.50
16:CP:80:PHE:H	16:CP:80:PHE:HD1	1.57	0.50
19:CS:77:THR:OG1	19:CS:78:ARG:HD2	2.12	0.50
25:CY:136:ALA:O	25:CY:139:LYS:HB2	2.12	0.50
27:D1:32:LYS:O	27:D1:33:LYS:O	2.30	0.50
32:D6:32:ASN:CG	32:D6:33:LYS:N	2.64	0.50
35:DA:1186:G:H2'	35:DA:1187:G:C5'	2.41	0.50
35:DA:1267:U:H2'	35:DA:1267:U:O2	2.10	0.50
35:DA:1486:A:N6	35:DA:1504:C:H42	2.08	0.50
35:DA:1590:U:H2'	35:DA:1591:G:C5'	2.30	0.50
35:DA:182:A:H2	35:DA:433:C:H1'	1.77	0.50
35:DA:528:A:C2	35:DA:2043:C:H5'	2.30	0.50
35:DA:2732:G:C3'	35:DA:2733:A:C5'	2.90	0.50
35:DA:2820:A:H8	39:DE:191:PRO:CB	2.24	0.50
35:DA:2828:C:O2'	35:DA:2829:C:H5'	2.12	0.50
35:DA:845:G:HO2'	35:DA:846:C:H5	1.56	0.50
36:DB:60:C:O2'	36:DB:61:G:H5'	2.12	0.50
38:DD:39:LYS:HB2	38:DD:62:TYR:HB2	1.94	0.50
38:DD:94:LEU:N	38:DD:94:LEU:CD1	2.72	0.50
41:DG:114:ILE:CG1	41:DG:117:PHE:HB2	2.42	0.50
42:DH:97:ARG:O	42:DH:125:VAL:HG21	2.12	0.50
35:DA:528:A:OP2	44:DN:114:ARG:HD2	2.12	0.50
44:DN:34:LEU:HD11	44:DN:116:LEU:O	2.11	0.50
46:DP:126:VAL:HG22	46:DP:145:PRO:HB3	1.94	0.50
46:DP:16:ARG:HD3	46:DP:18:ARG:N	2.21	0.50
47:DQ:25:ASP:CB	47:DQ:67:ARG:HH22	2.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:DR:9:LYS:HZ1	48:DR:42:LYS:HB3	1.76	0.50
49:DS:24:LEU:HD22	49:DS:24:LEU:H	1.77	0.50
50:DT:122:ASP:O	50:DT:126:ALA:HB2	2.12	0.50
50:DT:3:ARG:O	50:DT:5:ALA:N	2.45	0.50
51:DU:15:LYS:O	51:DU:19:LYS:HG3	2.12	0.50
51:DU:78:THR:O	51:DU:80:ILE:N	2.44	0.50
52:DV:8:GLY:C	52:DV:10:LYS:H	2.15	0.50
52:DV:96:ILE:HG23	52:DV:97:LYS:H	1.76	0.50
53:DW:18:ARG:HG2	53:DW:18:ARG:HH11	1.75	0.50
56:DZ:24:LEU:HD21	56:DZ:86:VAL:HG21	1.93	0.50
1:AA:1030(A):G:H2'	1:AA:1030(C):G:OP2	2.11	0.50
1:AA:1428:A:H61	1:AA:1473:A:N6	2.10	0.50
1:AA:1518:A:C2	1:AA:1519:A:C2	2.99	0.50
1:AA:582:U:H2'	1:AA:583:A:C8	2.46	0.50
1:AA:789:U:O2	1:AA:789:U:H2'	2.12	0.50
2:AB:87:ARG:O	2:AB:88:ALA:HB2	2.12	0.50
3:AC:40:ARG:HH11	3:AC:40:ARG:HG3	1.77	0.50
3:AC:69:HIS:N	3:AC:69:HIS:CD2	2.79	0.50
4:AD:153:ARG:CB	4:AD:153:ARG:HH11	2.24	0.50
7:AG:115:ARG:HB2	7:AG:118:VAL:HG21	1.93	0.50
7:AG:121:ALA:N	7:AG:124:LEU:HD12	2.27	0.50
7:AG:126:ASP:O	7:AG:130:GLY:N	2.34	0.50
11:AK:23:ALA:HA	11:AK:28:THR:CG2	2.42	0.50
13:AM:69:GLU:CB	13:AM:72:ALA:HB3	2.41	0.50
13:AM:70:LEU:O	13:AM:74:VAL:HG23	2.11	0.50
1:AA:1226:C:OP1	13:AM:96:LEU:HD13	2.10	0.50
15:AO:78:TYR:C	15:AO:80:ALA:N	2.65	0.50
16:AP:65:GLN:N	16:AP:65:GLN:OE1	2.45	0.50
1:AA:267:C:OP1	17:AQ:67:LYS:HB2	2.10	0.50
20:AT:53:LEU:O	20:AT:54:LYS:C	2.50	0.50
25:AY:15:GLN:OE1	25:AY:168:PHE:HZ	1.95	0.50
25:AY:37:LEU:CD1	25:AY:38:LEU:HG	2.42	0.50
26:B0:24:LYS:NZ	35:BA:2355:C:O2'	2.45	0.50
26:B0:29:GLN:HB2	26:B0:67:VAL:CG2	2.41	0.50
27:B1:47:GLN:NE2	27:B1:47:GLN:O	2.44	0.50
27:B1:83:GLU:C	27:B1:85:LEU:H	2.13	0.50
35:BA:1221:C:H2'	35:BA:1221(A):C:C6	2.45	0.50
35:BA:1681:G:O2'	35:BA:1762:A:H2'	2.12	0.50
35:BA:2030:A:H5''	35:BA:2031:A:OP1	2.10	0.50
35:BA:2313:C:H5'	35:BA:2313:C:H6	1.75	0.50
35:BA:2841:C:C2	35:BA:2877:G:C2	3.00	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:363(F):A:O2'	35:BA:364:C:H5	1.94	0.50
35:BA:419:C:C2	35:BA:420:C:C6	2.99	0.50
35:BA:442:G:O4'	40:BF:46:ARG:HG2	2.11	0.50
35:BA:451:C:H41	35:BA:453:C:H3'	1.77	0.50
35:BA:830:G:C4	35:BA:2448:A:C5	3.00	0.50
36:BB:4:C:H2'	36:BB:5:C:O4'	2.12	0.50
36:BB:56:G:H4'	36:BB:57:A:H8	1.76	0.50
36:BB:78:A:O2'	36:BB:79:C:H5'	2.12	0.50
38:BD:106:ILE:O	38:BD:106:ILE:HG23	2.11	0.50
38:BD:111:LEU:HD22	38:BD:115:GLN:OE1	2.11	0.50
35:BA:729:G:N7	38:BD:208:LYS:HB2	2.25	0.50
40:BF:25:PRO:HG3	40:BF:119:ARG:HA	1.93	0.50
40:BF:132:VAL:CG2	40:BF:133:ASN:H	2.04	0.50
40:BF:132:VAL:HG13	40:BF:133:ASN:N	2.26	0.50
40:BF:170:LEU:HD23	40:BF:173:VAL:CG2	2.42	0.50
40:BF:185:ASP:O	40:BF:189:THR:HG23	2.11	0.50
40:BF:89:VAL:HG12	40:BF:90:PHE:H	1.77	0.50
41:BG:132:ASN:ND2	41:BG:158:ALA:HB1	2.27	0.50
42:BH:18:GLU:CB	42:BH:25:LYS:HD2	2.42	0.50
44:BN:34:LEU:HD11	44:BN:116:LEU:O	2.12	0.50
44:BN:126:PRO:O	44:BN:127:ASP:HB2	2.12	0.50
44:BN:97:ARG:O	44:BN:98:VAL:C	2.50	0.50
49:BS:89:ARG:NE	49:BS:89:ARG:CA	2.73	0.50
50:BT:23:ARG:HB2	50:BT:24:PRO:HD2	1.93	0.50
50:BT:50:ILE:H	50:BT:50:ILE:HD12	1.76	0.50
51:BU:95:LEU:O	51:BU:98:LEU:HG	2.11	0.50
52:BV:60:GLU:HA	52:BV:100:ARG:O	2.12	0.50
52:BV:72:VAL:HG13	52:BV:88:ARG:HH22	1.76	0.50
53:BW:76:VAL:HB	53:BW:103:ILE:HG22	1.93	0.50
54:BX:46:ALA:C	54:BX:47:PHE:CD1	2.85	0.50
54:BX:47:PHE:CD1	54:BX:47:PHE:N	2.80	0.50
54:BX:56:THR:O	54:BX:57:LEU:HG	2.11	0.50
55:BY:28:LYS:CE	55:BY:37:VAL:HA	2.42	0.50
55:BY:96:ILE:HG13	55:BY:99:CYS:HB2	1.93	0.50
56:BZ:136:PHE:HD1	56:BZ:137:ILE:N	2.10	0.50
56:BZ:137:ILE:O	56:BZ:137:ILE:HG23	2.12	0.50
1:CA:1420:C:H2'	1:CA:1421:G:H8	1.76	0.50
1:CA:33:A:H2'	1:CA:34:C:C6	2.47	0.50
1:CA:516:U:O2'	1:CA:517:G:H5'	2.12	0.50
1:CA:640:A:H4'	8:CH:116:LYS:HZ1	1.76	0.50
1:CA:668:G:O2'	1:CA:669:U:H5'	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:154:ASN:CA	4:CD:159:ARG:HH21	2.24	0.50
8:CH:128:GLY:O	8:CH:129:VAL:HG13	2.12	0.50
8:CH:6:ILE:O	8:CH:10:LEU:HG	2.11	0.50
14:CN:33:VAL:HG12	14:CN:34:TYR:N	2.27	0.50
15:CO:16:ALA:HB1	15:CO:21:ASP:HB3	1.94	0.50
16:CP:65:GLN:N	16:CP:65:GLN:OE1	2.45	0.50
18:CR:72:ARG:O	18:CR:75:ILE:N	2.44	0.50
32:D6:30:THR:O	32:D6:31:PRO:C	2.50	0.50
32:D6:40:CYS:HB2	32:D6:46:HIS:CE1	2.43	0.50
34:D8:34:TRP:HZ3	34:D8:41:ILE:HG23	1.76	0.50
35:DA:1251:C:OP2	35:DA:1251:C:H2'	2.12	0.50
35:DA:1628:G:O2'	35:DA:1629:U:H5'	2.12	0.50
35:DA:1937:A:N7	35:DA:1939:U:H2'	2.27	0.50
35:DA:199:A:C6	35:DA:2434:A:C6	3.00	0.50
35:DA:2009:G:C6	35:DA:2010:G:N7	2.80	0.50
35:DA:2246:G:H2'	35:DA:2247:A:C8	2.47	0.50
35:DA:2313:C:H5'	35:DA:2313:C:H6	1.77	0.50
35:DA:2447:G:O6	35:DA:2504:U:O4	2.30	0.50
35:DA:2460:U:H2'	35:DA:2461:C:O4'	2.11	0.50
35:DA:2536:G:C6	35:DA:2537:U:N3	2.80	0.50
35:DA:260:G:N2	35:DA:261:G:H1'	2.26	0.50
35:DA:2678:C:C2	35:DA:2679:A:C8	3.00	0.50
35:DA:2683:C:H5''	50:DT:53:ARG:HH22	1.77	0.50
35:DA:304:G:H1	35:DA:313:C:N4	2.10	0.50
35:DA:304:G:H2'	35:DA:305:U:C6	2.47	0.50
33:D7:37:LYS:HG3	35:DA:458:G:C8	2.46	0.50
35:DA:501:A:H2'	35:DA:502:A:C8	2.47	0.50
35:DA:52:A:C2'	35:DA:53:A:H5'	2.42	0.50
33:D7:11:LYS:HE2	35:DA:686:G:H5''	1.94	0.50
35:DA:758:C:O2'	35:DA:759:G:H5'	2.12	0.50
35:DA:822:U:H2'	35:DA:822:U:O2	2.12	0.50
39:DE:52:LEU:HD12	39:DE:53:PRO:CD	2.40	0.50
40:DF:80:ALA:HB3	40:DF:83:PHE:CD1	2.47	0.50
41:DG:181:ARG:O	41:DG:182:LYS:C	2.50	0.50
42:DH:92:ILE:C	42:DH:94:TYR:H	2.15	0.50
43:DI:9:LEU:H	43:DI:13:GLY:HA2	1.76	0.50
44:DN:46:VAL:CG1	44:DN:47:ALA:N	2.64	0.50
35:DA:631:A:C4'	46:DP:65:ARG:HA	2.42	0.50
49:DS:95:HIS:O	49:DS:96:GLY:C	2.49	0.50
50:DT:108:ARG:NH1	50:DT:108:ARG:HB3	2.27	0.50
50:DT:53:ARG:CG	50:DT:53:ARG:HH11	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:DU:65:ILE:HD12	51:DU:65:ILE:N	2.26	0.50
55:DY:13:VAL:HG12	55:DY:14:LEU:N	2.27	0.50
56:DZ:56:VAL:C	56:DZ:57:ILE:HD12	2.32	0.50
1:AA:1230:C:O2'	1:AA:1231:G:H5'	2.11	0.50
1:AA:1260:C:H4'	1:AA:1284:C:H5'	1.93	0.50
1:AA:1512:U:H3	1:AA:1523:G:H1	1.60	0.50
1:AA:527:G:O2'	1:AA:528:C:H5'	2.12	0.50
1:AA:528:C:O2'	1:AA:529:G:H5'	2.12	0.50
1:AA:894:G:H2'	1:AA:895:G:C8	2.47	0.50
2:AB:215:LEU:O	2:AB:219:VAL:HG23	2.12	0.50
2:AB:80:ILE:HG13	2:AB:81:VAL:HG23	1.93	0.50
2:AB:73:THR:HB	2:AB:94:ASN:O	2.12	0.50
3:AC:150:LYS:HA	3:AC:169:ALA:HB2	1.94	0.50
4:AD:153:ARG:NH2	4:AD:181:MET:HG2	2.26	0.50
5:AE:105:VAL:HB	5:AE:106:PRO:CD	2.41	0.50
5:AE:102:ALA:HB1	5:AE:106:PRO:HG3	1.93	0.50
6:AF:7:ASN:C	6:AF:8:ILE:HG13	2.31	0.50
8:AH:40:ALA:HB2	8:AH:45:ILE:HD11	1.94	0.50
10:AJ:39:PRO:HB3	10:AJ:70:ARG:NH1	2.26	0.50
11:AK:28:THR:CG2	11:AK:29:ILE:N	2.74	0.50
11:AK:48:ILE:O	11:AK:50:TYR:N	2.41	0.50
11:AK:92:GLU:O	11:AK:95:ILE:HG12	2.10	0.50
12:AL:91:LYS:HG3	12:AL:91:LYS:O	2.12	0.50
19:AS:43:GLU:C	19:AS:45:VAL:N	2.63	0.50
1:AA:1226:C:OP1	19:AS:81:ARG:NH1	2.45	0.50
27:B1:92:LYS:N	27:B1:93:GLU:OE2	2.45	0.50
29:B3:19:GLN:C	29:B3:21:ALA:N	2.64	0.50
33:B7:18:PHE:O	33:B7:21:ARG:N	2.45	0.50
35:BA:1027:A:O2'	35:BA:1028:A:H5'	2.12	0.50
35:BA:1131:G:H1'	35:BA:1132:A:C8	2.47	0.50
35:BA:1695:G:H2'	35:BA:1696:G:C5'	2.42	0.50
35:BA:2491:U:C5'	35:BA:2570:G:H5''	2.24	0.50
35:BA:271(H):G:O2'	35:BA:271(I):G:H8	1.94	0.50
35:BA:2847:U:OP1	50:BT:98:LYS:HD3	2.12	0.50
35:BA:2854:G:H1	35:BA:2863:C:H42	1.59	0.50
35:BA:29:U:O2'	35:BA:30:G:H5'	2.11	0.50
35:BA:579:G:H2'	35:BA:580:C:C6	2.46	0.50
35:BA:614:U:O4'	35:BA:614:U:O2	2.30	0.50
35:BA:70:G:H21	35:BA:71:A:H62	1.60	0.50
36:BB:40:U:C4	36:BB:43:C:H5''	2.47	0.50
36:BB:75:G:N7	36:BB:76:G:N7	2.59	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:108:PRO:HB3	38:BD:143:HIS:CE1	2.47	0.50
40:BF:110:LEU:HD21	40:BF:181:LEU:CD2	2.41	0.50
40:BF:17:ARG:HG3	40:BF:17:ARG:HH11	1.77	0.50
41:BG:82:LEU:CD2	41:BG:86:MET:HB2	2.41	0.50
43:BI:112:LYS:H	43:BI:114:LEU:HG	1.77	0.50
45:BO:26:LYS:O	45:BO:27:GLY:O	2.30	0.50
35:BA:1246:A:P	46:BP:16:ARG:HH22	2.35	0.50
46:BP:90:ARG:HH11	46:BP:91:PHE:HB3	1.75	0.50
47:BQ:55:VAL:CG2	47:BQ:56:ARG:N	2.75	0.50
50:BT:106:SER:CA	50:BT:110:ILE:HD13	2.32	0.50
52:BV:3:ALA:O	52:BV:14:VAL:N	2.45	0.50
54:BX:7:VAL:O	54:BX:31:HIS:N	2.43	0.50
54:BX:35:THR:HG23	54:BX:36:LYS:H	1.76	0.50
55:BY:54:LYS:O	55:BY:55:TYR:O	2.29	0.50
56:BZ:112:ARG:O	56:BZ:113:ALA:CB	2.60	0.50
1:CA:1065:U:O2'	1:CA:1066:C:OP2	2.25	0.50
1:CA:1340:A:OP1	23:CW:36:A:H4'	2.11	0.50
1:CA:376:G:H2'	1:CA:377:G:H8	1.77	0.50
1:CA:525:C:H2'	1:CA:526:C:H6	1.76	0.50
1:CA:769:G:O2'	1:CA:770:C:H5'	2.11	0.50
1:CA:575:G:N1	1:CA:821:G:C5	2.80	0.50
1:CA:836:G:C6	1:CA:851:G:C6	3.00	0.50
1:CA:885:G:H2'	1:CA:886:G:H8	1.76	0.50
2:CB:81:VAL:HG22	2:CB:215:LEU:CG	2.42	0.50
4:CD:119:GLN:NE2	4:CD:123:HIS:NE2	2.60	0.50
4:CD:138:TYR:HD2	4:CD:139:ARG:N	2.09	0.50
4:CD:177:ASP:O	4:CD:180:GLY:N	2.43	0.50
4:CD:62:GLN:HB3	4:CD:66:ARG:HH22	1.77	0.50
5:CE:149:GLU:C	5:CE:151:LEU:H	2.14	0.50
7:CG:119:ARG:O	7:CG:122:HIS:N	2.45	0.50
5:CE:78:HIS:CD2	8:CH:104:ARG:HE	2.29	0.50
1:CA:1148:U:OP1	9:CI:7:THR:HG21	2.11	0.50
10:CJ:33:GLN:O	10:CJ:75:ILE:HG12	2.11	0.50
11:CK:109:VAL:HG13	18:CR:85:LEU:O	2.11	0.50
12:CL:104:VAL:O	12:CL:105:TYR:HB2	2.11	0.50
12:CL:55:VAL:O	12:CL:56:ALA:HB2	2.12	0.50
12:CL:60:LEU:CD2	12:CL:64:TYR:O	2.60	0.50
15:CO:78:TYR:O	15:CO:82:ILE:HG22	2.11	0.50
15:CO:82:ILE:HG12	15:CO:87:ILE:HB	1.93	0.50
16:CP:8:ARG:NH1	16:CP:8:ARG:HG2	2.24	0.50
1:CA:103:C:OP2	20:CT:14:LYS:HD2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:D0:77:ARG:HH22	35:DA:857:C:C5'	2.24	0.50
28:D2:41:ILE:C	28:D2:43:GLN:N	2.65	0.50
35:DA:1331:A:H2'	35:DA:1333:C:C5	2.46	0.50
35:DA:1338:G:H5'	35:DA:1339:G:OP2	2.10	0.50
35:DA:1382:G:O2'	35:DA:1383:C:H5'	2.11	0.50
35:DA:1407:C:H2'	35:DA:1407:C:O2	2.12	0.50
35:DA:141:A:H8	35:DA:1408:C:O2'	1.91	0.50
27:D1:41:ARG:NH1	35:DA:189:G:P	2.85	0.50
35:DA:1938:A:N1	35:DA:2590:A:H1'	2.27	0.50
35:DA:1982:C:OP1	35:DA:1982:C:H3'	2.12	0.50
35:DA:2417:C:H2'	35:DA:2418:A:H8	1.77	0.50
35:DA:2605:U:H2'	35:DA:2606:C:C6	2.47	0.50
35:DA:2701:C:C3'	35:DA:2702:U:H5''	2.29	0.50
35:DA:2718:G:H2'	35:DA:2719:G:H8	1.76	0.50
35:DA:2880:C:O3'	48:DR:90:ARG:NH1	2.44	0.50
35:DA:644:A:O2'	35:DA:645:C:H5''	2.12	0.50
35:DA:679:C:O2	35:DA:680:G:C8	2.65	0.50
35:DA:71:A:H4'	35:DA:72:U:C5'	2.41	0.50
28:D2:41:ILE:HD13	35:DA:94(A):G:O2'	2.11	0.50
38:DD:73:VAL:HG13	38:DD:120:GLY:CA	2.42	0.50
38:DD:24:ILE:HD13	38:DD:24:ILE:O	2.12	0.50
38:DD:260:ARG:CZ	38:DD:264:LYS:HD3	2.42	0.50
41:DG:132:ASN:CG	41:DG:158:ALA:HA	2.31	0.50
43:DI:123:LEU:HG	43:DI:142:VAL:HB	1.93	0.50
43:DI:79:ILE:CG2	43:DI:81:VAL:HG23	2.42	0.50
43:DI:82:ARG:HH11	43:DI:82:ARG:HG3	1.77	0.50
44:DN:3:THR:C	44:DN:5:VAL:H	2.14	0.50
46:DP:21:ARG:NH1	46:DP:21:ARG:HG3	2.26	0.50
47:DQ:77:LYS:O	47:DQ:79:LEU:N	2.38	0.50
48:DR:52:ILE:CD1	48:DR:79:LEU:HD21	2.42	0.50
48:DR:56:LYS:HE2	48:DR:94:TYR:CE2	2.47	0.50
49:DS:26:LEU:O	49:DS:26:LEU:CD2	2.60	0.50
49:DS:74:ALA:CB	49:DS:103:GLU:HG3	2.41	0.50
52:DV:60:GLU:HA	52:DV:100:ARG:O	2.12	0.50
54:DX:32:PRO:HD3	54:DX:72:LYS:HZ2	1.76	0.50
55:DY:2:ARG:N	55:DY:5:MET:CE	2.75	0.50
56:DZ:93:ASP:OD1	56:DZ:94:GLU:HG2	2.12	0.50
1:AA:1323:G:H2'	1:AA:1324:A:C8	2.47	0.50
1:AA:1333:A:H3'	1:AA:1334:G:H8	1.77	0.50
1:AA:19:C:P	5:AE:127:ASN:HD22	2.35	0.50
2:AB:164:VAL:HB	2:AB:186:ALA:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:103:ASN:O	4:AD:106:TYR:N	2.45	0.50
5:AE:41:VAL:HG13	5:AE:113:ALA:HA	1.93	0.50
8:AH:28:ALA:CB	8:AH:57:PRO:O	2.60	0.50
17:AQ:45:HIS:HB2	17:AQ:69:LYS:CE	2.42	0.50
20:AT:26:ASN:ND2	20:AT:26:ASN:H	2.10	0.50
25:AY:16:LYS:O	25:AY:19:GLU:HB2	2.11	0.50
26:B0:51:VAL:HG13	26:B0:60:PHE:O	2.12	0.50
27:B1:53:VAL:HG13	27:B1:54:ALA:N	2.26	0.50
27:B1:61:ARG:O	27:B1:62:VAL:O	2.30	0.50
27:B1:85:LEU:C	27:B1:87:PRO:CD	2.69	0.50
35:BA:1141:U:C5'	35:BA:1142(A):A:O4'	2.60	0.50
35:BA:1485:G:H2'	35:BA:1486:A:C8	2.46	0.50
35:BA:1747:G:O2'	35:BA:1747(A):G:H5'	2.12	0.50
35:BA:2090:G:H2'	35:BA:2091:U:O4'	2.11	0.50
35:BA:261:G:C2	35:BA:262:A:C8	3.00	0.50
35:BA:2678:C:C2	35:BA:2679:A:C8	3.00	0.50
35:BA:2881:C:H2'	35:BA:2882:A:C8	2.45	0.50
35:BA:90:U:O2'	35:BA:92:A:H5''	2.12	0.50
35:BA:995:C:C2	51:BU:57:PHE:CE2	2.99	0.50
35:BA:1658:C:OP1	39:BE:132:HIS:O	2.29	0.50
39:BE:125:GLY:N	39:BE:135:HIS:O	2.44	0.50
40:BF:28:ILE:HD13	40:BF:28:ILE:H	1.74	0.50
41:BG:167:GLU:O	41:BG:171:ALA:N	2.41	0.50
43:BI:92:VAL:CG1	43:BI:120:ILE:HD12	2.32	0.50
45:BO:63:VAL:HB	45:BO:102:VAL:HG12	1.94	0.50
45:BO:62:VAL:CG1	45:BO:65:THR:HG22	2.42	0.50
47:BQ:51:ARG:C	47:BQ:54:MET:HB3	2.31	0.50
47:BQ:9:TYR:C	47:BQ:9:TYR:CD2	2.84	0.50
48:BR:118:GLU:HA	48:BR:118:GLU:OE1	2.12	0.50
48:BR:41:ALA:O	48:BR:44:LEU:N	2.44	0.50
51:BU:78:THR:C	51:BU:80:ILE:N	2.65	0.50
52:BV:19:LYS:CE	52:BV:19:LYS:HA	2.40	0.50
52:BV:23:GLU:O	52:BV:24:LYS:O	2.30	0.50
53:BW:29:LEU:HD21	53:BW:33:ARG:CZ	2.40	0.50
53:BW:31:GLU:O	53:BW:32:ALA:C	2.50	0.50
54:BX:33:LYS:C	54:BX:35:THR:H	2.14	0.50
54:BX:57:LEU:CB	54:BX:76:ARG:HD2	2.39	0.50
56:BZ:29:TYR:HE2	56:BZ:87:ASP:HB3	1.77	0.50
56:BZ:61:LEU:HB2	56:BZ:65:GLN:CB	2.40	0.50
1:CA:1305:G:H5''	21:CU:5:ASP:N	2.27	0.50
1:CA:1379:G:C4	1:CA:1380:U:H5	2.30	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1423:G:H2'	1:CA:1424:C:H6	1.77	0.50
1:CA:1499:A:H2'	1:CA:1500:A:C8	2.43	0.50
1:CA:183:G:H1	1:CA:194:C:H42	1.59	0.50
1:CA:286:G:H2'	1:CA:287:U:C6	2.47	0.50
1:CA:408:A:C6	1:CA:409:G:C5	3.00	0.50
1:CA:709:G:H2'	1:CA:710:G:H8	1.77	0.50
2:CB:224:GLN:HG3	2:CB:229:VAL:CG2	2.42	0.50
2:CB:22:LYS:HZ3	2:CB:22:LYS:HA	1.72	0.50
2:CB:69:LEU:HD13	2:CB:91:PRO:O	2.12	0.50
4:CD:192:GLU:OE2	4:CD:192:GLU:N	2.45	0.50
4:CD:24:GLU:O	4:CD:27:TYR:HB2	2.12	0.50
6:CF:42:GLU:C	6:CF:44:GLY:N	2.65	0.50
6:CF:67:MET:HE2	6:CF:72:VAL:H	1.77	0.50
11:CK:22:HIS:HB3	11:CK:29:ILE:CG1	2.35	0.50
11:CK:57:THR:O	11:CK:60:ALA:HB3	2.12	0.50
13:CM:108:ARG:H	13:CM:108:ARG:CD	2.12	0.50
16:CP:40:ASP:HB3	16:CP:48:TRP:HB3	1.93	0.50
17:CQ:10:VAL:HG12	17:CQ:53:LEU:HD12	1.94	0.50
17:CQ:31:LEU:HG	17:CQ:32:TYR:CD2	2.47	0.50
22:CV:32:U:O5'	22:CV:32:U:H6	1.95	0.50
25:CY:143:LEU:C	25:CY:143:LEU:HD23	2.32	0.50
25:CY:32:ARG:C	25:CY:103:ILE:HD13	2.32	0.50
25:CY:38:LEU:HA	25:CY:41:LEU:HD11	1.92	0.50
26:D0:42:GLY:HA3	35:DA:2331:G:O4'	2.11	0.50
27:D1:34:THR:HG23	27:D1:34:THR:O	2.10	0.50
31:D5:41:PRO:HG2	31:D5:44:THR:HG1	1.76	0.50
33:D7:18:PHE:O	33:D7:19:ARG:C	2.49	0.50
34:D8:30:ARG:HH21	46:DP:62:LEU:CB	2.20	0.50
34:D8:35:GLN:NE2	34:D8:36:LYS:HZ2	2.10	0.50
35:DA:1195:G:H2'	35:DA:1196:C:C6	2.47	0.50
35:DA:1578:U:O2	35:DA:1578:U:H2'	2.11	0.50
35:DA:1643:G:H2'	35:DA:1644:C:C6	2.42	0.50
35:DA:1682:G:H2'	35:DA:1683:C:C6	2.46	0.50
35:DA:1691:C:O2'	35:DA:1692:U:H5'	2.12	0.50
35:DA:1904:G:N2	35:DA:1905:C:H1'	2.27	0.50
35:DA:1926:U:H2'	35:DA:1928:A:OP2	2.11	0.50
25:CY:130:ARG:HH21	35:DA:1942:C:C2'	2.25	0.50
35:DA:2010:G:H5''	53:DW:42:ARG:HB2	1.93	0.50
35:DA:2284:C:H42	35:DA:2384:G:H1	1.59	0.50
35:DA:2637:U:H1'	35:DA:2782:G:N2	2.27	0.50
35:DA:2892:A:C4	35:DA:2893:G:H1'	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:363(A):A:C2	35:DA:363(B):G:C8	3.00	0.50
38:DD:125:ILE:HG22	38:DD:125:ILE:O	2.11	0.50
38:DD:183:ARG:HG2	38:DD:183:ARG:NH1	2.27	0.50
38:DD:222:ARG:O	38:DD:224:ALA:N	2.44	0.50
38:DD:34:VAL:CG1	38:DD:34:VAL:O	2.60	0.50
38:DD:92:ILE:CA	38:DD:107:ALA:HB2	2.41	0.50
39:DE:38:THR:CG2	39:DE:39:PRO:HD2	2.41	0.50
40:DF:197:ASP:OD1	40:DF:198:ALA:N	2.44	0.50
40:DF:30:PRO:O	40:DF:33:LEU:HB3	2.11	0.50
40:DF:41:LEU:HA	40:DF:44:ARG:HD3	1.93	0.50
42:DH:125:VAL:O	42:DH:125:VAL:HG12	2.11	0.50
42:DH:85:LYS:HB3	42:DH:133:VAL:O	2.12	0.50
43:DI:36:ALA:O	43:DI:37:VAL:HG23	2.12	0.50
44:DN:103:VAL:O	44:DN:106:MET:HB2	2.12	0.50
44:DN:113:GLY:O	44:DN:114:ARG:C	2.51	0.50
44:DN:34:LEU:CD2	44:DN:120:LEU:HD23	2.41	0.50
44:DN:33:LEU:HD23	44:DN:38:HIS:CE1	2.47	0.50
45:DO:6:THR:CG2	45:DO:7:TYR:N	2.69	0.50
46:DP:101:VAL:HA	46:DP:107:LYS:H	1.77	0.50
46:DP:84:ASN:CG	46:DP:116:GLY:HA3	2.31	0.50
40:DF:34:TRP:CZ3	46:DP:12:ALA:HA	2.47	0.50
48:DR:4:LEU:C	48:DR:6:SER:H	2.16	0.50
49:DS:46:VAL:HG12	49:DS:47:THR:N	2.26	0.50
49:DS:85:VAL:CG2	49:DS:86:ALA:H	2.16	0.50
50:DT:45:PHE:CZ	50:DT:74:ARG:HB2	2.46	0.50
51:DU:55:ARG:HA	51:DU:58:ARG:CG	2.42	0.50
51:DU:66:ASN:ND2	51:DU:70:ARG:NH2	2.39	0.50
52:DV:34:GLU:HB3	52:DV:62:LEU:CD1	2.37	0.50
52:DV:2:PHE:O	52:DV:3:ALA:HB2	2.12	0.50
52:DV:52:VAL:C	52:DV:54:GLY:N	2.65	0.50
52:DV:72:VAL:HG13	52:DV:88:ARG:NH2	2.27	0.50
52:DV:72:VAL:HA	52:DV:88:ARG:HH22	1.77	0.50
53:DW:18:ARG:HG2	53:DW:18:ARG:NH1	2.27	0.50
28:D2:26:ARG:NE	54:DX:5:TYR:HB3	2.27	0.50
56:DZ:14:LYS:C	56:DZ:16:SER:H	2.15	0.50
56:DZ:76:LEU:N	56:DZ:84:GLU:HB2	2.25	0.50
1:AA:1234:C:H4'	1:AA:1364:U:H1'	1.93	0.49
1:AA:603:U:H2'	1:AA:604:G:H8	1.76	0.49
1:AA:640:A:C2'	1:AA:641:U:H5'	2.42	0.49
6:AF:75:LEU:HD23	6:AF:79:LEU:HD21	1.94	0.49
9:AI:45:ALA:O	9:AI:49:PRO:HD2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:119:LYS:C	12:AL:120:TYR:CD1	2.85	0.49
12:AL:85:ILE:CD1	12:AL:98:TYR:HB3	2.42	0.49
15:AO:74:ASP:OD2	15:AO:77:ARG:HG2	2.12	0.49
1:AA:237:C:C4'	17:AQ:25:ARG:HH12	2.18	0.49
17:AQ:57:VAL:HA	17:AQ:77:VAL:HG23	1.94	0.49
18:AR:43:PHE:H	18:AR:43:PHE:HD1	1.58	0.49
19:AS:50:ALA:HB1	19:AS:57:HIS:HB3	1.94	0.49
25:AY:156:ARG:HH22	47:BQ:80:GLU:CB	2.25	0.49
25:AY:29:ARG:HB3	25:AY:32:ARG:NE	2.26	0.49
25:AY:51:PRO:HB2	25:AY:53:ASN:ND2	2.27	0.49
26:B0:37:LEU:C	26:B0:38:VAL:CG2	2.81	0.49
26:B0:29:GLN:O	26:B0:66:VAL:HA	2.12	0.49
31:B5:20:ARG:O	31:B5:23:HIS:CD2	2.64	0.49
33:B7:25:PRO:HA	33:B7:28:ARG:CZ	2.42	0.49
35:BA:106:C:H1'	55:BY:2:ARG:CZ	2.42	0.49
35:BA:1141:U:OP2	44:BN:63:THR:HG21	2.11	0.49
35:BA:1863:G:H1	35:BA:1879:C:N4	2.09	0.49
35:BA:2282:G:C4	35:BA:2425:A:N6	2.80	0.49
35:BA:271(R):G:O2'	35:BA:271(S):G:H5'	2.12	0.49
35:BA:2732:G:C3'	35:BA:2733:A:C5'	2.88	0.49
35:BA:2777:G:H4'	35:BA:2778:A:H5'	1.94	0.49
35:BA:295:G:C2	35:BA:296:C:C6	3.00	0.49
35:BA:359:A:H2'	35:BA:360:G:C8	2.47	0.49
35:BA:71:A:H8	35:BA:71:A:H5'	1.77	0.49
35:BA:778:G:C6	35:BA:779:U:N3	2.80	0.49
35:BA:833:U:H2'	35:BA:834:C:H6	1.75	0.49
35:BA:927:G:O6	35:BA:928:G:C2	2.64	0.49
35:BA:848:G:C2	35:BA:933:A:H1'	2.47	0.49
35:BA:939:G:O2'	35:BA:940:G:H5'	2.12	0.49
38:BD:211:ARG:NH1	38:BD:211:ARG:HG2	2.26	0.49
38:BD:63:ARG:CZ	38:BD:86:PRO:HD2	2.42	0.49
39:BE:115:GLY:HA2	39:BE:157:ALA:CB	2.42	0.49
39:BE:59:VAL:HG13	39:BE:60:ASN:N	2.25	0.49
40:BF:21:ALA:C	40:BF:23:ASP:H	2.16	0.49
40:BF:65:TRP:HH2	40:BF:75:HIS:HD2	1.60	0.49
41:BG:116:ASP:CG	41:BG:117:PHE:N	2.65	0.49
42:BH:92:ILE:C	42:BH:94:TYR:H	2.15	0.49
35:BA:833:U:P	46:BP:45:LEU:HD11	2.52	0.49
35:BA:2406:U:C4	46:BP:72:PRO:HD2	2.47	0.49
49:BS:74:ALA:CB	49:BS:103:GLU:HG3	2.42	0.49
49:BS:18:ILE:C	49:BS:20:ARG:H	2.15	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:88:ILE:HG22	50:BT:89:VAL:H	1.76	0.49
51:BU:15:LYS:O	51:BU:19:LYS:HG3	2.11	0.49
54:BX:53:LYS:NZ	54:BX:55:ASN:ND2	2.60	0.49
56:BZ:144:LEU:N	56:BZ:144:LEU:HD22	2.26	0.49
56:BZ:33:LEU:HD12	56:BZ:34:ASN:N	2.26	0.49
1:CA:1312:G:H1	1:CA:1325:C:H42	1.60	0.49
1:CA:23:C:O2'	1:CA:24:U:H5'	2.12	0.49
1:CA:25:C:O2'	1:CA:26:A:H5'	2.12	0.49
1:CA:338:A:H2'	1:CA:339:C:C6	2.47	0.49
1:CA:375:U:OP1	16:CP:69:THR:HG21	2.12	0.49
1:CA:676:A:O2'	1:CA:677:U:H5'	2.12	0.49
1:CA:903:G:H2'	1:CA:904:C:C6	2.46	0.49
5:CE:126:ARG:O	5:CE:127:ASN:C	2.48	0.49
5:CE:43:LEU:CD2	5:CE:132:ALA:HB1	2.33	0.49
5:CE:80:ILE:HD11	5:CE:91:LEU:HD22	1.94	0.49
6:CF:88:VAL:CG1	6:CF:89:MET:N	2.75	0.49
7:CG:29:LYS:HG3	7:CG:101:LEU:HB3	1.93	0.49
7:CG:15:ASP:HB2	7:CG:23:VAL:HB	1.94	0.49
8:CH:85:ARG:HA	8:CH:135:CYS:HB3	1.94	0.49
9:CI:28:VAL:CG1	9:CI:29:ASN:N	2.75	0.49
10:CJ:39:PRO:HB3	10:CJ:70:ARG:NH1	2.27	0.49
13:CM:11:ARG:O	13:CM:13:LYS:HG3	2.12	0.49
13:CM:89:GLY:O	13:CM:92:HIS:HB2	2.12	0.49
14:CN:57:ARG:HG2	14:CN:58:LYS:N	2.27	0.49
20:CT:26:ASN:ND2	20:CT:26:ASN:H	2.09	0.49
26:D0:11:ARG:O	26:D0:11:ARG:HG2	2.12	0.49
34:D8:58:ILE:C	34:D8:61:LEU:HG	2.32	0.49
35:DA:1467:C:H42	35:DA:1525:G:H1	1.60	0.49
35:DA:1885:A:H3'	35:DA:1886:C:C6	2.47	0.49
35:DA:1948:G:H2'	35:DA:1949:G:H8	1.77	0.49
35:DA:1964:G:H3'	35:DA:1965:C:C5'	2.42	0.49
35:DA:2025:C:H2'	35:DA:2026:C:C6	2.47	0.49
35:DA:2069:G:C2	35:DA:2070:G:C8	3.00	0.49
35:DA:2115:G:O2'	35:DA:2116:G:H5''	2.12	0.49
35:DA:2119:A:C3'	35:DA:2120:G:C5'	2.90	0.49
35:DA:2119:A:C2'	35:DA:2120:G:H5''	2.42	0.49
35:DA:2189:U:H2'	35:DA:2190:G:O4'	2.13	0.49
35:DA:2678:C:O2	35:DA:2678:C:H2'	2.12	0.49
35:DA:2714:G:O2'	35:DA:2715:C:H5'	2.11	0.49
35:DA:271(R):G:O2'	35:DA:271(S):G:H5'	2.12	0.49
35:DA:510:C:O2'	35:DA:511:U:H5'	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:820:A:O2'	35:DA:821:A:H5'	2.12	0.49
35:DA:860:U:O2	35:DA:860:U:O4'	2.29	0.49
38:DD:113:VAL:O	38:DD:115:GLN:N	2.35	0.49
38:DD:152:GLY:O	38:DD:154:LYS:HG3	2.12	0.49
39:DE:11:MET:HB2	39:DE:23:VAL:O	2.11	0.49
40:DF:46:ARG:HA	40:DF:46:ARG:NH1	2.22	0.49
42:DH:113:VAL:HG21	42:DH:151:ILE:HG21	1.94	0.49
42:DH:163:TYR:N	42:DH:163:TYR:CD1	2.80	0.49
43:DI:98:ALA:HB1	43:DI:109:ILE:CG1	2.42	0.49
44:DN:31:ALA:O	44:DN:34:LEU:N	2.45	0.49
44:DN:97:ARG:O	44:DN:101:HIS:N	2.40	0.49
46:DP:71:VAL:HG22	46:DP:72:PRO:HG3	1.94	0.49
48:DR:75:LEU:O	48:DR:79:LEU:HB2	2.12	0.49
35:DA:1011:G:OP2	51:DU:70:ARG:NH2	2.45	0.49
55:DY:76:CYS:HB3	55:DY:96:ILE:HD11	1.93	0.49
56:DZ:115:GLY:CA	56:DZ:175:VAL:O	2.59	0.49
1:AA:1464:G:O2'	1:AA:1465:C:H5'	2.12	0.49
1:AA:236:G:H2'	1:AA:237:C:C6	2.48	0.49
1:AA:542:G:C6	1:AA:543:C:N4	2.81	0.49
1:AA:556:C:C2'	1:AA:557:G:H5'	2.42	0.49
1:AA:661:G:H1	1:AA:744:C:H42	1.59	0.49
1:AA:833:U:H2'	1:AA:834:C:C6	2.47	0.49
1:AA:940:C:H2'	1:AA:941:G:H8	1.77	0.49
2:AB:74:LYS:HG3	2:AB:74:LYS:O	2.11	0.49
3:AC:35:GLU:HA	3:AC:38:ARG:HG2	1.93	0.49
3:AC:47:LEU:HG	3:AC:52:LEU:HD22	1.95	0.49
8:AH:119:LEU:HD12	8:AH:123:GLU:C	2.32	0.49
10:AJ:16:LEU:HA	10:AJ:19:SER:OG	2.12	0.49
12:AL:66:VAL:HG21	12:AL:98:TYR:CD1	2.47	0.49
15:AO:56:LEU:O	15:AO:57:LEU:C	2.50	0.49
16:AP:39:TYR:OH	16:AP:41:PRO:HB3	2.13	0.49
17:AQ:47:PRO:HG2	17:AQ:48:GLU:OE1	2.12	0.49
18:AR:37:VAL:O	18:AR:38:GLU:C	2.50	0.49
25:AY:171:LYS:HA	25:AY:174:GLN:HE22	1.75	0.49
27:B1:83:GLU:C	27:B1:85:LEU:N	2.65	0.49
29:B3:54:VAL:CG1	29:B3:55:ARG:N	2.75	0.49
31:B5:21:SER:O	31:B5:23:HIS:N	2.45	0.49
34:B8:39:LYS:O	34:B8:42:ARG:HB3	2.12	0.49
35:BA:1324:G:H3'	35:BA:1325:G:C4'	2.42	0.49
35:BA:1428:C:H41	35:BA:1569:A:H3'	1.76	0.49
35:BA:1615:C:H5	35:BA:1617:C:C2	2.30	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2247:A:O2'	35:BA:2248:C:H5'	2.12	0.49
35:BA:2417:C:H2'	35:BA:2418:A:H8	1.77	0.49
35:BA:2589:A:H2'	35:BA:2590:A:C8	2.46	0.49
35:BA:336:C:H4'	55:BY:7:VAL:HG13	1.94	0.49
35:BA:658:C:H2'	35:BA:659:C:C6	2.47	0.49
39:BE:105:THR:HB	39:BE:197:ILE:HG12	1.93	0.49
41:BG:169:ALA:O	41:BG:170:ARG:C	2.49	0.49
41:BG:80:PHE:O	41:BG:81:LYS:O	2.29	0.49
42:BH:113:VAL:HG21	42:BH:151:ILE:HG21	1.94	0.49
44:BN:27:ALA:O	44:BN:28:THR:C	2.50	0.49
46:BP:79:ARG:O	46:BP:111:ARG:HB2	2.11	0.49
47:BQ:88:GLY:C	47:BQ:90:VAL:H	2.14	0.49
49:BS:99:LYS:O	49:BS:100:ALA:C	2.50	0.49
50:BT:61:PHE:CE2	50:BT:76:PHE:HB3	2.47	0.49
35:BA:583:G:OP2	51:BU:10:ARG:NH1	2.45	0.49
53:BW:18:ARG:HG2	53:BW:18:ARG:NH1	2.26	0.49
54:BX:7:VAL:HB	54:BX:8:ILE:HD12	1.94	0.49
56:BZ:19:ARG:HH11	56:BZ:19:ARG:CB	2.25	0.49
1:CA:1030(A):G:H2'	1:CA:1030(C):G:OP2	2.12	0.49
1:CA:1405:G:H21	1:CA:1517:G:N2	2.09	0.49
1:CA:189(I):G:H2'	1:CA:189(J):G:C8	2.47	0.49
1:CA:337:C:O2'	1:CA:338:A:H5'	2.11	0.49
1:CA:429:U:H4'	1:CA:430:A:O5'	2.12	0.49
1:CA:707:C:H2'	1:CA:707:C:O2	2.12	0.49
1:CA:959:A:C2'	1:CA:960:U:H4'	2.42	0.49
2:CB:167:PRO:CD	2:CB:188:ALA:HB2	2.42	0.49
3:CC:47:LEU:HG	3:CC:52:LEU:HD22	1.94	0.49
4:CD:194:LEU:O	4:CD:195:ALA:HB3	2.12	0.49
5:CE:135:THR:O	5:CE:139:LEU:HG	2.13	0.49
6:CF:28:ARG:HG3	6:CF:28:ARG:NH1	2.26	0.49
6:CF:55:ASP:OD2	6:CF:86:ARG:NH2	2.45	0.49
8:CH:9:MET:O	8:CH:12:ARG:N	2.45	0.49
8:CH:83:ILE:HB	8:CH:137:VAL:CG1	2.38	0.49
11:CK:20:TYR:CE2	11:CK:83:ILE:HB	2.47	0.49
13:CM:83:ASP:OD1	13:CM:86:CYS:HB2	2.12	0.49
16:CP:20:VAL:CG2	16:CP:21:VAL:N	2.74	0.49
17:CQ:56:VAL:HG23	17:CQ:78:GLU:O	2.11	0.49
19:CS:12:ASP:CB	19:CS:15:LEU:HD23	2.41	0.49
27:D1:76:ARG:HE	27:D1:76:ARG:HA	1.76	0.49
27:D1:75:GLU:CB	27:D1:76:ARG:NH2	2.70	0.49
27:D1:94:LEU:CD1	27:D1:95:LEU:N	2.74	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D2:20:GLU:C	28:D2:22:GLU:H	2.15	0.49
29:D3:27:GLY:O	29:D3:35:ARG:HD2	2.12	0.49
30:D4:29:PRO:C	30:D4:31:ILE:N	2.65	0.49
31:D5:51:TYR:C	31:D5:56:LYS:HG2	2.32	0.49
35:DA:1233:C:H2'	35:DA:1234:U:H5'	1.93	0.49
35:DA:1269:A:H2'	35:DA:1270:C:C6	2.47	0.49
35:DA:769:G:H4'	35:DA:1379:A:N1	2.27	0.49
35:DA:1778:U:H5	35:DA:1784:A:C2	2.30	0.49
35:DA:1914:C:H3'	35:DA:1914:C:O2	2.12	0.49
35:DA:2606:C:C2'	35:DA:2607:G:H5'	2.42	0.49
35:DA:2732:G:O2'	35:DA:2733:A:H5'	2.11	0.49
35:DA:542:C:C2'	35:DA:543:C:OP1	2.59	0.49
35:DA:910:A:C4	47:DQ:13:GLN:OE1	2.65	0.49
38:DD:133:LEU:O	38:DD:136:ILE:HG13	2.12	0.49
38:DD:226:MET:CB	38:DD:230:ASP:HB2	2.40	0.49
38:DD:25:THR:O	38:DD:25:THR:HG23	2.10	0.49
35:DA:1816:G:N7	38:DD:62:TYR:CE1	2.80	0.49
39:DE:125:GLY:N	39:DE:135:HIS:O	2.45	0.49
39:DE:4:ILE:HD11	39:DE:28:ALA:HB1	1.94	0.49
40:DF:155:LEU:HD13	40:DF:174:VAL:HB	1.94	0.49
35:DA:2667:C:H1'	42:DH:109:PHE:CE2	2.47	0.49
47:DQ:101:ARG:HH11	47:DQ:101:ARG:HG3	1.76	0.49
50:DT:109:GLU:HA	50:DT:112:ARG:HD2	1.93	0.49
51:DU:28:ARG:HA	51:DU:34:LYS:HB3	1.94	0.49
51:DU:95:LEU:O	51:DU:98:LEU:HG	2.11	0.49
52:DV:61:VAL:HG23	52:DV:100:ARG:HG2	1.93	0.49
52:DV:63:GLY:O	52:DV:64:HIS:HB3	2.11	0.49
35:DA:1312:U:OP2	54:DX:62:LYS:HE2	2.13	0.49
55:DY:87:LYS:CG	55:DY:88:LYS:N	2.74	0.49
1:AA:1131:G:H2'	1:AA:1132:C:C5	2.46	0.49
1:AA:1421:G:H2'	1:AA:1422:G:H8	1.76	0.49
1:AA:376:G:H2'	1:AA:377:G:H8	1.76	0.49
1:AA:377:G:O2'	1:AA:378:G:H5'	2.11	0.49
2:AB:145:LEU:HD22	2:AB:149:LEU:HD12	1.94	0.49
2:AB:21:ARG:O	2:AB:23:ARG:N	2.45	0.49
4:AD:120:LEU:HA	4:AD:125:HIS:CD2	2.47	0.49
4:AD:150:GLU:HG2	4:AD:151:LYS:H	1.78	0.49
4:AD:199:ASN:OD1	4:AD:201:GLN:HB2	2.12	0.49
4:AD:68:TYR:N	4:AD:68:TYR:CD1	2.80	0.49
6:AF:71:ARG:O	6:AF:72:VAL:C	2.49	0.49
7:AG:106:GLN:HA	7:AG:109:ASN:HD22	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:92:SER:OG	7:AG:94:ARG:NE	2.45	0.49
11:AK:20:TYR:CE2	11:AK:83:ILE:HB	2.47	0.49
11:AK:87:THR:HG22	11:AK:88:GLY:N	2.28	0.49
11:AK:58:PRO:HA	11:AK:90:GLY:CA	2.43	0.49
1:AA:502:G:OP1	12:AL:118:SER:N	2.43	0.49
13:AM:64:TRP:NE1	13:AM:66:LEU:HD12	2.26	0.49
14:AN:57:ARG:HG2	14:AN:58:LYS:N	2.27	0.49
1:AA:472:A:H4'	16:AP:82:GLN:HE22	1.78	0.49
17:AQ:85:VAL:O	17:AQ:89:LEU:HB2	2.12	0.49
25:AY:156:ARG:NH2	26:B0:6:ALA:HB2	2.27	0.49
28:B2:58:ALA:O	28:B2:60:LEU:N	2.46	0.49
33:B7:8:ASN:ND2	33:B7:9:ARG:N	2.58	0.49
35:BA:1246:A:P	46:BP:16:ARG:HH12	2.35	0.49
35:BA:1903:G:C2	35:BA:1904:G:C8	3.00	0.49
35:BA:2009:G:H1'	48:BR:107:ASP:C	2.32	0.49
35:BA:2018:G:C6	35:BA:2019:A:C5	3.00	0.49
35:BA:2030:A:H4'	35:BA:2031:A:C8	2.47	0.49
35:BA:2049:G:O2'	35:BA:2050:C:H5'	2.12	0.49
35:BA:2242:G:H2'	35:BA:2243:U:O4'	2.12	0.49
35:BA:2282:G:O2'	35:BA:2283:C:OP2	2.25	0.49
35:BA:309:G:H4'	55:BY:18:GLY:HA3	1.93	0.49
35:BA:559:G:N2	51:BU:49:HIS:HD2	2.09	0.49
35:BA:733:G:H8	35:BA:733:G:O5'	1.96	0.49
35:BA:685:A:C5	35:BA:774:A:C2	3.00	0.49
35:BA:892:G:H3'	35:BA:892:G:N3	2.28	0.49
36:BB:103:G:H2'	36:BB:104:U:C6	2.47	0.49
39:BE:161:GLY:O	39:BE:163:GLU:HG2	2.12	0.49
40:BF:114:VAL:HG21	40:BF:202:PHE:CE2	2.47	0.49
40:BF:9:ILE:HG12	40:BF:14:PRO:CA	2.42	0.49
42:BH:89:ILE:HD13	42:BH:90:LYS:N	2.24	0.49
43:BI:25:TYR:HE2	43:BI:29:TYR:CD2	2.30	0.49
43:BI:55:ALA:O	43:BI:59:ALA:HB2	2.12	0.49
45:BO:86:ILE:O	45:BO:93:PRO:HA	2.13	0.49
46:BP:106:LEU:CD1	46:BP:112:LEU:HB2	2.42	0.49
47:BQ:141:GLN:OXT	56:BZ:98:MET:HE1	2.12	0.49
48:BR:116:LEU:O	48:BR:117:VAL:CB	2.60	0.49
49:BS:92:TYR:C	49:BS:92:TYR:CD1	2.86	0.49
50:BT:62:THR:CG2	50:BT:75:ILE:HA	2.23	0.49
1:CA:1132:C:O2'	1:CA:1133:G:H5'	2.11	0.49
1:CA:1152:A:H2'	1:CA:1153:C:C6	2.46	0.49
1:CA:1234:C:C4'	1:CA:1364:U:H1'	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1251:A:O2'	1:CA:1252:A:H5'	2.13	0.49
1:CA:1480:G:N2	1:CA:1481:U:H1'	2.28	0.49
1:CA:243:A:O2'	1:CA:244:U:OP2	2.29	0.49
1:CA:666:G:H2'	1:CA:667:G:H8	1.77	0.49
1:CA:781:A:C2'	1:CA:782:A:H5'	2.41	0.49
2:CB:101:MET:O	2:CB:105:PHE:CA	2.58	0.49
2:CB:50:GLU:O	2:CB:51:LEU:C	2.51	0.49
2:CB:74:LYS:O	2:CB:74:LYS:HG3	2.13	0.49
3:CC:91:LEU:HB2	3:CC:99:VAL:HG21	1.93	0.49
9:CI:118:LYS:H	9:CI:121:ARG:HB3	1.77	0.49
11:CK:127:LYS:O	11:CK:129:SER:N	2.44	0.49
13:CM:69:GLU:HA	13:CM:70:LEU:N	2.27	0.49
16:CP:60:LEU:C	16:CP:62:VAL:N	2.66	0.49
19:CS:16:LEU:C	19:CS:20:LEU:HG	2.31	0.49
25:CY:84:ARG:C	25:CY:86:SER:H	2.15	0.49
28:D2:25:VAL:HA	28:D2:28:LYS:HB2	1.95	0.49
29:D3:26:LEU:O	29:D3:27:GLY:C	2.51	0.49
31:D5:21:SER:O	31:D5:23:HIS:N	2.46	0.49
33:D7:15:THR:HG22	33:D7:16:HIS:NE2	2.27	0.49
35:DA:1195:G:H2'	35:DA:1196:C:H6	1.76	0.49
35:DA:1317:A:C2	35:DA:1318:C:C2	3.00	0.49
35:DA:1435:G:H2'	35:DA:1436:G:O4'	2.12	0.49
35:DA:1490:A:H5'	35:DA:1491:G:OP2	2.11	0.49
35:DA:16:G:H2'	35:DA:17:G:C8	2.48	0.49
35:DA:1922:G:O2'	35:DA:1923:U:H5'	2.12	0.49
35:DA:201:C:O2'	35:DA:202:U:H5'	2.12	0.49
35:DA:1136:G:N3	35:DA:2038:G:H4'	2.28	0.49
35:DA:2379:G:H2'	35:DA:2380:C:C6	2.47	0.49
35:DA:2259:G:H1'	35:DA:2427:C:O2	2.11	0.49
35:DA:2463:C:O2'	35:DA:2464:C:H5'	2.12	0.49
35:DA:2519:U:H3'	35:DA:2519:U:OP1	2.12	0.49
35:DA:2847:U:OP1	50:DT:98:LYS:HD3	2.12	0.49
35:DA:335:C:H4'	55:DY:73:ARG:CZ	2.42	0.49
35:DA:336:C:H4'	55:DY:7:VAL:HG13	1.93	0.49
35:DA:754:C:H2'	35:DA:755:C:C6	2.47	0.49
35:DA:876:C:H2'	35:DA:877:U:C6	2.47	0.49
37:DC:40:THR:HG21	37:DC:215:THR:CB	2.42	0.49
35:DA:2203:U:H4'	38:DD:151:LYS:HE3	1.94	0.49
40:DF:152:GLU:OE1	40:DF:191:ARG:HD2	2.12	0.49
41:DG:38:VAL:CB	41:DG:158:ALA:HB3	2.42	0.49
42:DH:105:LEU:CD2	42:DH:105:LEU:N	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:124:GLU:HB2	42:DH:132:ARG:O	2.13	0.49
42:DH:156:ALA:O	42:DH:158:HIS:N	2.36	0.49
43:DI:55:ALA:O	43:DI:59:ALA:HB2	2.12	0.49
43:DI:79:ILE:CB	43:DI:81:VAL:HG23	2.42	0.49
44:DN:126:PRO:O	44:DN:127:ASP:HB2	2.11	0.49
44:DN:26:LEU:CD2	44:DN:99:LEU:HD11	2.40	0.49
44:DN:84:LYS:O	44:DN:85:ILE:HD13	2.12	0.49
45:DO:43:VAL:C	45:DO:45:GLU:H	2.15	0.49
35:DA:1246:A:P	46:DP:16:ARG:HH22	2.35	0.49
46:DP:5:ASP:OD2	46:DP:6:LEU:HD22	2.12	0.49
47:DQ:34:LEU:HD22	47:DQ:121:ALA:HB3	1.94	0.49
49:DS:18:ILE:C	49:DS:20:ARG:H	2.15	0.49
49:DS:34:HIS:CE1	49:DS:55:ALA:HB2	2.45	0.49
49:DS:86:ALA:C	49:DS:106:ARG:HG3	2.33	0.49
51:DU:105:VAL:HG12	51:DU:109:LEU:HD11	1.92	0.49
51:DU:86:ALA:HB2	51:DU:116:ALA:CB	2.42	0.49
53:DW:59:VAL:O	53:DW:61:ASN:N	2.45	0.49
54:DX:56:THR:O	54:DX:57:LEU:HG	2.12	0.49
56:DZ:33:LEU:HD11	56:DZ:35:ARG:CB	2.42	0.49
56:DZ:33:LEU:HD11	56:DZ:35:ARG:HB2	1.93	0.49
35:DA:1040:C:H5''	56:DZ:46:LYS:NZ	2.26	0.49
56:DZ:47:VAL:HG12	56:DZ:51:ALA:CB	2.42	0.49
1:AA:1307:U:H2'	1:AA:1308:U:H6	1.74	0.49
1:AA:1431:C:H42	1:AA:1469:G:H1	1.60	0.49
1:AA:1480:G:C2	1:AA:1481:U:C2	3.00	0.49
2:AB:69:LEU:HB2	2:AB:159:PRO:CG	2.42	0.49
3:AC:8:ILE:O	3:AC:10:PHE:N	2.45	0.49
4:AD:14:ARG:C	4:AD:16:GLY:H	2.16	0.49
8:AH:21:LYS:CG	8:AH:22:GLU:N	2.75	0.49
8:AH:80:ILE:O	8:AH:80:ILE:HG22	2.12	0.49
10:AJ:55:LYS:O	10:AJ:56:HIS:CG	2.65	0.49
11:AK:127:LYS:O	11:AK:129:SER:N	2.42	0.49
13:AM:37:THR:O	13:AM:39:ILE:HG13	2.12	0.49
13:AM:89:GLY:O	13:AM:92:HIS:HB2	2.13	0.49
17:AQ:29:HIS:CA	17:AQ:36:ILE:HD11	2.42	0.49
17:AQ:47:PRO:HG2	17:AQ:48:GLU:OE2	2.12	0.49
25:AY:29:ARG:HG3	25:AY:114:LEU:HD13	1.94	0.49
27:B1:18:ILE:HG12	27:B1:43:TYR:HD1	1.78	0.49
28:B2:32:LEU:CG	28:B2:33:MET:N	2.75	0.49
35:BA:1496:A:H8	35:BA:1577:C:O2'	1.95	0.49
35:BA:1819:A:H1'	35:BA:1821:A:C5	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:202:U:O2'	35:BA:203:C:H5'	2.11	0.49
35:BA:203:C:H2'	35:BA:204:A:H8	1.78	0.49
35:BA:2354:G:O2'	35:BA:2355:C:H5'	2.12	0.49
35:BA:234:C:H2'	35:BA:235:U:C6	2.46	0.49
35:BA:2728:U:H2'	35:BA:2728:U:O2	2.12	0.49
35:BA:2814:C:H2'	35:BA:2815:C:C6	2.42	0.49
35:BA:576:U:OP1	35:BA:2503:A:OP1	2.30	0.49
37:BC:65:PRO:HG2	37:BC:189:ILE:CB	2.43	0.49
39:BE:52:LEU:HB2	39:BE:76:ARG:CB	2.36	0.49
41:BG:110:ALA:HB1	41:BG:140:ILE:HD13	1.93	0.49
41:BG:174:GLU:CA	41:BG:178:PHE:HB2	2.41	0.49
42:BH:130:ARG:HB2	42:BH:130:ARG:HH11	1.76	0.49
42:BH:18:GLU:HB3	42:BH:25:LYS:HD2	1.95	0.49
44:BN:26:LEU:HG	44:BN:30:ILE:CD1	2.42	0.49
44:BN:61:ARG:HG3	44:BN:61:ARG:HH11	1.77	0.49
44:BN:96:GLU:O	44:BN:97:ARG:C	2.51	0.49
46:BP:47:ASP:CB	46:BP:51:PHE:HB2	2.33	0.49
52:BV:88:ARG:HG3	52:BV:88:ARG:NH1	2.28	0.49
52:BV:8:GLY:C	52:BV:10:LYS:H	2.14	0.49
54:BX:82:GLN:HG3	54:BX:83:VAL:H	1.72	0.49
1:CA:1029:C:H1'	1:CA:1033:G:N1	2.28	0.49
1:CA:1494:G:C2	1:CA:1495:U:C6	3.01	0.49
1:CA:405:U:H5''	1:CA:406:G:O4'	2.11	0.49
1:CA:564:C:C5	17:CQ:31:LEU:HD11	2.48	0.49
1:CA:640:A:C2'	1:CA:641:U:H5'	2.42	0.49
2:CB:24:TRP:O	2:CB:25:ASN:HB2	2.13	0.49
6:CF:10:LEU:HD21	6:CF:22:GLU:HB3	1.94	0.49
8:CH:11:THR:CA	8:CH:14:ARG:HH12	2.25	0.49
10:CJ:16:LEU:O	10:CJ:16:LEU:HD13	2.12	0.49
13:CM:117:VAL:CG1	13:CM:118:ALA:N	2.76	0.49
13:CM:97:PRO:O	13:CM:98:VAL:N	2.45	0.49
14:CN:29:ARG:HH11	14:CN:29:ARG:HG3	1.78	0.49
18:CR:37:VAL:CG2	18:CR:38:GLU:H	2.05	0.49
18:CR:43:PHE:H	18:CR:43:PHE:HD1	1.54	0.49
25:CY:140:LEU:HD21	25:CY:157:ALA:HB1	1.94	0.49
25:CY:32:ARG:CB	25:CY:103:ILE:HD13	2.42	0.49
25:CY:84:ARG:HE	25:CY:92:PRO:HD2	1.77	0.49
29:D3:54:VAL:CG1	29:D3:55:ARG:N	2.75	0.49
34:D8:2:PRO:C	34:D8:4:MET:H	2.15	0.49
34:D8:6:THR:O	34:D8:8:LYS:N	2.46	0.49
35:DA:1190:G:O3'	46:DP:35:HIS:HB3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2019:A:N6	35:DA:2020:A:C6	2.81	0.49
35:DA:2206:G:H21	35:DA:2207:G:C5'	2.21	0.49
35:DA:2319:G:OP2	35:DA:2319:G:H4'	2.12	0.49
35:DA:2666:C:O2'	35:DA:2667:C:H5'	2.13	0.49
35:DA:2675:A:OP1	45:DO:31:LYS:HB2	2.12	0.49
35:DA:776:G:H4'	35:DA:777:A:O5'	2.11	0.49
35:DA:875:G:HO2'	56:DZ:151:HIS:CD2	2.29	0.49
36:DB:103:G:H2'	36:DB:104:U:C6	2.47	0.49
37:DC:83:ILE:HG22	37:DC:83:ILE:O	2.12	0.49
40:DF:167:ALA:HB1	40:DF:173:VAL:CG1	2.42	0.49
40:DF:45:ARG:HG2	40:DF:97:TYR:CG	2.47	0.49
41:DG:7:LEU:O	41:DG:11:TYR:N	2.45	0.49
44:DN:79:PRO:HG2	44:DN:80:GLY:H	1.77	0.49
45:DO:35:VAL:H	45:DO:65:THR:HG21	1.77	0.49
46:DP:47:ASP:HB2	46:DP:51:PHE:CD2	2.48	0.49
47:DQ:120:ILE:HA	47:DQ:123:HIS:HD2	1.77	0.49
35:DA:1276:A:H1'	48:DR:16:HIS:HE1	1.76	0.49
48:DR:41:ALA:C	48:DR:43:GLU:N	2.65	0.49
49:DS:72:ALA:O	49:DS:73:LEU:C	2.50	0.49
51:DU:112:ARG:HG2	51:DU:112:ARG:HH11	1.77	0.49
35:DA:992:C:O3'	52:DV:75:PHE:CE2	2.65	0.49
52:DV:96:ILE:HG23	52:DV:97:LYS:N	2.26	0.49
53:DW:37:ARG:HG3	53:DW:37:ARG:NH1	2.25	0.49
1:AA:1042:G:O2'	1:AA:1043:C:H5'	2.11	0.49
1:AA:1287:A:H2	1:AA:1353:G:N3	2.10	0.49
1:AA:498:U:H2'	1:AA:499:A:C5'	2.41	0.49
1:AA:59:A:H5'	1:AA:60:A:H5'	1.94	0.49
1:AA:788:U:C2	1:AA:789:U:C6	3.00	0.49
2:AB:157:ARG:HG2	2:AB:158:LEU:N	2.27	0.49
2:AB:165:VAL:O	2:AB:187:LEU:O	2.31	0.49
2:AB:187:LEU:HB2	2:AB:201:ILE:HB	1.95	0.49
3:AC:172:ARG:HB3	3:AC:174:PRO:HD3	1.93	0.49
4:AD:80:GLU:HB3	4:AD:84:LYS:HZ1	1.77	0.49
5:AE:37:ARG:HH11	5:AE:37:ARG:HG2	1.77	0.49
8:AH:86:ILE:CG1	8:AH:135:CYS:HA	2.39	0.49
9:AI:17:VAL:CG1	9:AI:81:ILE:HD13	2.43	0.49
11:AK:88:GLY:N	11:AK:91:ARG:HB2	2.27	0.49
12:AL:104:VAL:O	12:AL:105:TYR:HB2	2.12	0.49
12:AL:119:LYS:H	12:AL:119:LYS:HD2	1.77	0.49
15:AO:43:LEU:C	15:AO:45:VAL:H	2.16	0.49
16:AP:17:TYR:CD1	16:AP:17:TYR:N	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:82:MET:HA	17:AQ:85:VAL:CG2	2.43	0.49
18:AR:76:LEU:HD22	18:AR:76:LEU:N	2.27	0.49
32:B6:28:ARG:HG2	32:B6:28:ARG:HH11	1.77	0.49
32:B6:37:ARG:O	32:B6:48:VAL:O	2.30	0.49
34:B8:21:LYS:O	34:B8:23:VAL:HG23	2.12	0.49
35:BA:1227:G:O2'	35:BA:1228:G:H5'	2.12	0.49
35:BA:1410:G:C6	35:BA:1411:C:N4	2.80	0.49
35:BA:1608:A:O3'	35:BA:1609:A:H3'	2.13	0.49
35:BA:1613:G:C6	35:BA:1619:G:C6	3.00	0.49
35:BA:2545:G:N2	35:BA:2546:U:H1'	2.27	0.49
35:BA:2689:U:H4'	35:BA:2690:C:H6	1.77	0.49
35:BA:2704:C:H2'	35:BA:2705:A:O4'	2.13	0.49
35:BA:2762:G:H2'	35:BA:2763:G:C5'	2.42	0.49
35:BA:543:C:C4	35:BA:551:G:N1	2.81	0.49
35:BA:542:C:C2'	35:BA:543:C:OP1	2.60	0.49
35:BA:671:C:H5	46:BP:36:LYS:NZ	2.11	0.49
35:BA:795:C:H2'	35:BA:796:C:H6	1.74	0.49
35:BA:820:A:O2'	35:BA:821:A:H5'	2.12	0.49
35:BA:871:U:H4'	47:BQ:69:PHE:CE2	2.47	0.49
36:BB:60:C:C2	36:BB:61:G:C8	3.00	0.49
36:BB:65:C:O2'	36:BB:66:A:H5'	2.12	0.49
38:BD:25:THR:CG2	38:BD:82:ILE:N	2.70	0.49
39:BE:6:GLY:CA	39:BE:27:LEU:O	2.60	0.49
40:BF:53:THR:O	40:BF:54:ARG:C	2.49	0.49
41:BG:55:LYS:O	41:BG:57:ALA:N	2.45	0.49
44:BN:26:LEU:HD21	44:BN:30:ILE:HD11	1.94	0.49
45:BO:107:ARG:NH1	50:BT:36:GLU:N	2.60	0.49
47:BQ:101:ARG:HG3	47:BQ:101:ARG:HH11	1.76	0.49
48:BR:101:ALA:O	48:BR:102:GLU:HB2	2.12	0.49
48:BR:49:ASP:O	48:BR:52:ILE:N	2.46	0.49
48:BR:56:LYS:NZ	48:BR:88:ARG:O	2.43	0.49
35:BA:2838:G:N2	48:BR:93:GLY:HA3	2.27	0.49
49:BS:107:GLU:HG3	49:BS:108:GLY:N	2.27	0.49
49:BS:15:ARG:C	49:BS:17:ARG:N	2.65	0.49
49:BS:34:HIS:CD2	49:BS:53:SER:CB	2.94	0.49
53:BW:20:VAL:HG21	53:BW:47:VAL:HG21	1.94	0.49
55:BY:18:GLY:C	55:BY:20:TYR:H	2.15	0.49
35:BA:336:C:H4'	55:BY:7:VAL:CG1	2.42	0.49
1:CA:1405:G:H1'	1:CA:1518:A:O2'	2.12	0.49
1:CA:1469:G:H2'	1:CA:1470:G:C8	2.47	0.49
1:CA:156:G:C2	1:CA:166:G:N1	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:184:G:O4'	1:CA:224:C:H4'	2.13	0.49
1:CA:865:A:H2	1:CA:918:A:H4'	1.77	0.49
4:CD:18:LYS:CE	4:CD:31:CYS:SG	3.01	0.49
4:CD:87:GLY:C	4:CD:89:THR:H	2.16	0.49
7:CG:133:GLY:O	7:CG:136:LYS:HB2	2.13	0.49
7:CG:47:CYS:HB3	7:CG:58:PRO:HG3	1.95	0.49
8:CH:122:ARG:NH1	8:CH:122:ARG:HB2	2.27	0.49
1:CA:875:C:O2'	8:CH:14:ARG:NH1	2.46	0.49
8:CH:80:ILE:HG22	8:CH:80:ILE:O	2.12	0.49
17:CQ:56:VAL:CG2	17:CQ:78:GLU:HG3	2.42	0.49
25:CY:67:VAL:HB	25:CY:98:ALA:HB1	1.93	0.49
29:D3:12:PRO:O	29:D3:14:GLY:N	2.45	0.49
35:DA:1428:C:H41	35:DA:1569:A:H3'	1.76	0.49
35:DA:1434:A:O2'	35:DA:1435:G:H5'	2.13	0.49
35:DA:1467:C:N4	35:DA:1525:G:H1	2.10	0.49
35:DA:1795:C:H1'	35:DA:1901:A:OP1	2.12	0.49
35:DA:1854:A:H3'	35:DA:1855:G:C8	2.46	0.49
35:DA:2014:A:C2	35:DA:2015:A:N1	2.81	0.49
35:DA:2119:A:C3'	35:DA:2120:G:H5''	2.43	0.49
35:DA:2408:U:C2	35:DA:2409:G:C8	3.00	0.49
35:DA:2547:U:H2'	35:DA:2548:G:H8	1.77	0.49
35:DA:2620:C:OP1	39:DE:152:LYS:O	2.30	0.49
35:DA:2822:G:H2'	35:DA:2823:A:H5''	1.94	0.49
35:DA:686:G:N2	35:DA:788:A:N6	2.59	0.49
35:DA:737:C:C2'	35:DA:738:G:O5'	2.61	0.49
36:DB:83:G:C2	36:DB:84:C:C6	3.01	0.49
38:DD:176:ARG:NH1	38:DD:176:ARG:HG2	2.23	0.49
35:DA:1824:G:OP1	38:DD:52:ARG:HD3	2.12	0.49
38:DD:43:ARG:CB	38:DD:54:ARG:HB2	2.43	0.49
38:DD:95:LEU:N	38:DD:95:LEU:HD12	2.27	0.49
40:DF:125:LEU:HD12	40:DF:196:LEU:HD22	1.95	0.49
40:DF:1:MET:O	40:DF:2:LYS:C	2.50	0.49
41:DG:160:VAL:CG1	41:DG:161:THR:N	2.74	0.49
41:DG:32:PRO:O	41:DG:172:LEU:HD11	2.13	0.49
42:DH:113:VAL:HG21	42:DH:151:ILE:CG2	2.42	0.49
42:DH:89:ILE:CG1	42:DH:90:LYS:N	2.75	0.49
46:DP:47:ASP:CB	46:DP:51:PHE:HB2	2.34	0.49
47:DQ:110:THR:OG1	47:DQ:113:GLN:HG3	2.12	0.49
47:DQ:20:ALA:CB	47:DQ:99:PRO:O	2.60	0.49
50:DT:75:ILE:CD1	50:DT:75:ILE:N	2.76	0.49
51:DU:92:ARG:HG2	51:DU:92:ARG:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:DX:30:VAL:O	54:DX:31:HIS:C	2.50	0.49
56:DZ:71:VAL:HG13	56:DZ:88:PHE:CE2	2.48	0.49
1:AA:259:G:O2'	1:AA:260:G:H5'	2.13	0.49
1:AA:358:U:H2'	1:AA:359:U:C6	2.47	0.49
1:AA:390:C:O5'	1:AA:390:C:H6	1.96	0.49
1:AA:447:G:N2	1:AA:488:C:N4	2.59	0.49
1:AA:645:C:H2'	1:AA:646:U:C6	2.47	0.49
1:AA:788:U:H2'	1:AA:789:U:H6	1.78	0.49
2:AB:80:ILE:HG12	2:AB:211:ILE:HG21	1.95	0.49
3:AC:52:LEU:O	3:AC:53:ALA:HB2	2.11	0.49
7:AG:133:GLY:O	7:AG:136:LYS:HB2	2.13	0.49
8:AH:128:GLY:O	8:AH:129:VAL:HG13	2.12	0.49
8:AH:7:ALA:O	8:AH:11:THR:N	2.32	0.49
9:AI:6:GLY:HA3	9:AI:84:ALA:N	2.27	0.49
11:AK:27:ASN:OD1	11:AK:28:THR:N	2.45	0.49
11:AK:62:GLN:O	11:AK:64:ALA:N	2.45	0.49
12:AL:57:LYS:O	12:AL:58:VAL:HG23	2.13	0.49
16:AP:48:TRP:HE3	16:AP:49:LEU:HB3	1.76	0.49
19:AS:77:THR:OG1	19:AS:78:ARG:HD2	2.12	0.49
20:AT:26:ASN:HD22	20:AT:27:LYS:H	1.60	0.49
21:AU:18:TYR:HA	21:AU:22:ARG:HD3	1.94	0.49
23:AW:23:G:O2'	23:AW:24:C:H5''	2.11	0.49
23:AW:53:G:N2	23:AW:54:G:N7	2.61	0.49
25:AY:156:ARG:NH2	26:B0:6:ALA:CB	2.75	0.49
29:B3:36:VAL:HG23	29:B3:36:VAL:O	2.12	0.49
35:BA:1210:A:O2'	35:BA:1211:U:OP2	2.25	0.49
35:BA:1214:A:H2'	35:BA:1215:G:C8	2.47	0.49
35:BA:11:G:O2'	35:BA:12:U:H5'	2.13	0.49
35:BA:2206:G:H21	35:BA:2207:G:C5'	2.22	0.49
35:BA:271(V):G:C2	35:BA:271(W):G:H1'	2.48	0.49
35:BA:2822:G:OP2	39:BE:110:GLY:O	2.31	0.49
35:BA:2830:G:H5'	39:BE:58:ARG:CZ	2.43	0.49
35:BA:371:A:C4	35:BA:373:U:C4	3.01	0.49
35:BA:376:C:O2'	35:BA:377:C:H5'	2.12	0.49
35:BA:848:G:H2'	35:BA:849:A:C8	2.47	0.49
35:BA:957:A:H5'	47:BQ:76:LYS:HG3	1.93	0.49
37:BC:56:GLN:HB2	37:BC:168:THR:HA	1.94	0.49
38:BD:211:ARG:O	38:BD:214:TRP:N	2.46	0.49
39:BE:133:LYS:N	39:BE:134:ILE:HD13	2.27	0.49
39:BE:4:ILE:HD11	39:BE:28:ALA:HB1	1.94	0.49
39:BE:9:VAL:HG23	39:BE:9:VAL:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:45:ARG:HD2	40:BF:46:ARG:H	1.78	0.49
40:BF:62:ARG:NH2	40:BF:64:ILE:HA	2.28	0.49
41:BG:163:ALA:HB1	41:BG:168:GLU:HB2	1.95	0.49
41:BG:19:LEU:O	41:BG:23:PHE:N	2.45	0.49
43:BI:58:LEU:O	43:BI:58:LEU:HD23	2.12	0.49
35:BA:598:G:C5'	46:BP:15:ARG:HD2	2.37	0.49
47:BQ:32:TYR:HA	47:BQ:132:VAL:O	2.12	0.49
47:BQ:85:LYS:HG3	47:BQ:86:GLY:H	1.78	0.49
48:BR:44:LEU:HD13	48:BR:48:VAL:HG23	1.93	0.49
51:BU:40:PHE:N	51:BU:40:PHE:HD2	2.08	0.49
51:BU:66:ASN:ND2	51:BU:70:ARG:NH2	2.38	0.49
51:BU:76:TYR:O	51:BU:80:ILE:HG12	2.13	0.49
53:BW:26:GLY:H	53:BW:71:VAL:CG1	2.25	0.49
55:BY:31:LEU:HB2	55:BY:36:ALA:N	2.23	0.49
56:BZ:75:ASN:O	56:BZ:84:GLU:N	2.45	0.49
1:CA:1149:C:H2'	1:CA:1150:U:C6	2.43	0.49
1:CA:1342:C:H2'	1:CA:1343:G:C8	2.48	0.49
1:CA:1490:C:O2'	1:CA:1491:G:H5'	2.12	0.49
2:CB:164:VAL:HB	2:CB:186:ALA:HB2	1.95	0.49
3:CC:105:GLU:CG	3:CC:106:VAL:H	2.17	0.49
4:CD:65:ARG:HH11	4:CD:72:GLU:CA	2.25	0.49
6:CF:83:ASP:O	6:CF:85:VAL:N	2.44	0.49
8:CH:97:VAL:HG22	8:CH:98:LYS:N	2.28	0.49
11:CK:21:ILE:CD1	11:CK:21:ILE:N	2.70	0.49
11:CK:21:ILE:CD1	11:CK:82:VAL:HG13	2.37	0.49
12:CL:58:VAL:CG1	12:CL:59:ARG:N	2.76	0.49
19:CS:20:LEU:O	19:CS:23:ASN:HB3	2.12	0.49
20:CT:58:LYS:HE3	20:CT:62:LEU:CD1	2.42	0.49
24:CX:16:U:H2'	24:CX:17:U:C5	2.48	0.49
28:D2:44:LEU:C	28:D2:46:GLN:N	2.61	0.49
31:D5:32:PRO:O	31:D5:33:CYS:CB	2.60	0.49
33:D7:30:VAL:CG1	33:D7:33:ARG:HH22	2.25	0.49
35:DA:1827:C:H2'	35:DA:1828:G:O4'	2.13	0.49
35:DA:1853:A:N6	35:DA:1889:A:C8	2.80	0.49
35:DA:1786:A:C6	35:DA:1938:A:N7	2.80	0.49
35:DA:2321:G:N3	35:DA:2321:G:H2'	2.27	0.49
35:DA:2366:A:H2'	35:DA:2367:G:O4'	2.12	0.49
35:DA:2704:C:H2'	35:DA:2705:A:O4'	2.12	0.49
35:DA:2864:G:H2'	35:DA:2865:U:O4'	2.12	0.49
35:DA:445:C:H5''	51:DU:3:ARG:HB3	1.94	0.49
35:DA:836:G:C5	35:DA:837:C:C4	2.99	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:955:C:H5'	35:DA:956:G:OP2	2.13	0.49
36:DB:79:C:C2'	36:DB:80:U:H5'	2.42	0.49
38:DD:193:VAL:HG13	38:DD:193:VAL:O	2.12	0.49
38:DD:257:LEU:HD23	38:DD:258:LYS:O	2.12	0.49
38:DD:262:ARG:O	38:DD:264:LYS:N	2.46	0.49
39:DE:167:VAL:CG2	39:DE:168:MET:N	2.75	0.49
40:DF:21:ALA:O	40:DF:23:ASP:N	2.46	0.49
40:DF:47:GLY:HA3	40:DF:95:ARG:O	2.12	0.49
41:DG:28:VAL:O	41:DG:31:VAL:CG1	2.61	0.49
42:DH:41:MET:HE1	42:DH:55:PRO:HD2	1.95	0.49
43:DI:77:LEU:HD21	43:DI:101:LEU:HD13	1.95	0.49
44:DN:97:ARG:O	44:DN:98:VAL:C	2.49	0.49
46:DP:47:ASP:OD1	46:DP:50:ARG:HG3	2.12	0.49
47:DQ:134:ARG:C	47:DQ:136:ALA:H	2.16	0.49
48:DR:21:TYR:OH	48:DR:43:GLU:HG2	2.12	0.49
49:DS:17:ARG:HG3	49:DS:18:ILE:HD13	1.95	0.49
49:DS:90:GLY:C	49:DS:92:TYR:N	2.62	0.49
50:DT:58:ASN:ND2	50:DT:58:ASN:N	2.60	0.49
51:DU:24:TYR:HB2	51:DU:29:SER:OG	2.13	0.49
35:DA:1248:G:H2'	51:DU:2:PRO:O	2.12	0.49
53:DW:70:TYR:CE2	53:DW:108:GLY:HA3	2.45	0.49
53:DW:34:ASN:HA	53:DW:37:ARG:HB3	1.93	0.49
54:DX:52:VAL:H	54:DX:80:ILE:HG22	1.77	0.49
55:DY:14:LEU:HD12	55:DY:15:VAL:N	2.18	0.49
55:DY:45:VAL:HA	55:DY:62:GLU:CG	2.15	0.49
1:AA:1055:A:C2	1:AA:1056:U:H1'	2.48	0.49
1:AA:1293:G:O2'	1:AA:1294:G:H8	1.96	0.49
1:AA:482:A:H3'	1:AA:483:C:C6	2.48	0.49
1:AA:782:A:N6	1:AA:801:U:C6	2.81	0.49
2:AB:82:ARG:HA	2:AB:92:TYR:CE1	2.48	0.49
3:AC:107:GLN:N	3:AC:107:GLN:CD	2.65	0.49
3:AC:47:LEU:CD2	3:AC:52:LEU:HD13	2.43	0.49
4:AD:104:VAL:O	4:AD:107:ARG:HB2	2.12	0.49
4:AD:128:VAL:CG1	4:AD:129:ASN:N	2.72	0.49
4:AD:202:LEU:O	4:AD:203:VAL:C	2.51	0.49
4:AD:62:GLN:O	4:AD:63:LYS:C	2.51	0.49
4:AD:96:LEU:C	4:AD:98:GLU:N	2.65	0.49
5:AE:149:GLU:C	5:AE:151:LEU:N	2.65	0.49
5:AE:78:HIS:CD2	8:AH:104:ARG:NE	2.80	0.49
1:AA:1349:A:OP2	9:AI:118:LYS:NZ	2.45	0.49
3:AC:30:ARG:HD2	14:AN:38:GLY:HA3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:86:GLU:HB3	17:AQ:90:ILE:HG12	1.95	0.49
18:AR:63:GLN:OE1	18:AR:63:GLN:HA	2.12	0.49
25:AY:172:ALA:O	25:AY:173:ASP:C	2.50	0.49
28:B2:25:VAL:O	28:B2:29:LYS:HG3	2.12	0.49
28:B2:47:ASN:HA	28:B2:51:ARG:HB3	1.95	0.49
35:BA:1278:A:OP1	48:BR:36:THR:HG22	2.12	0.49
35:BA:1504:C:O2'	35:BA:1505:C:C5'	2.61	0.49
35:BA:151:C:H2'	35:BA:152:G:C8	2.48	0.49
35:BA:1639:U:H2'	35:BA:1640:C:C5'	2.40	0.49
35:BA:1788:C:C2	35:BA:1789:A:C8	3.01	0.49
35:BA:1791:A:OP2	35:BA:1791:A:H8	1.95	0.49
35:BA:1910:G:O2'	35:BA:1911:U:H5'	2.12	0.49
35:BA:2024:G:H2'	35:BA:2025:C:C6	2.47	0.49
35:BA:2246:G:N2	35:BA:2426:A:H1'	2.26	0.49
32:B6:27:LYS:HG3	35:BA:2286:A:OP2	2.11	0.49
35:BA:2650:U:H2'	35:BA:2651:C:C6	2.48	0.49
35:BA:2682:U:O2	39:BE:22:PRO:HB3	2.12	0.49
35:BA:2732:G:O2'	35:BA:2733:A:H5'	2.11	0.49
35:BA:442:G:C4'	40:BF:46:ARG:HD3	2.43	0.49
35:BA:859:G:H5'	35:BA:2268:A:O2'	2.12	0.49
35:BA:884:C:H4'	35:BA:892:G:N7	2.28	0.49
38:BD:92:ILE:C	38:BD:92:ILE:HD12	2.33	0.49
39:BE:103:ASP:CG	39:BE:201:THR:HA	2.33	0.49
40:BF:112:MET:O	40:BF:113:ALA:C	2.51	0.49
40:BF:31:HIS:HB2	46:BP:13:ASN:OD1	2.12	0.49
40:BF:7:TYR:HB3	40:BF:16:GLY:C	2.33	0.49
41:BG:37:VAL:HG13	41:BG:158:ALA:O	2.13	0.49
44:BN:54:VAL:HB	44:BN:122:VAL:HG22	1.94	0.49
44:BN:93:THR:C	44:BN:94:HIS:CG	2.85	0.49
45:BO:43:VAL:C	45:BO:45:GLU:H	2.16	0.49
45:BO:44:LYS:O	45:BO:45:GLU:HB3	2.12	0.49
46:BP:51:PHE:O	46:BP:52:GLU:HB2	2.12	0.49
47:BQ:134:ARG:NH1	56:BZ:119:GLU:OE2	2.45	0.49
47:BQ:50:ALA:O	47:BQ:54:MET:HB2	2.12	0.49
49:BS:66:ALA:O	49:BS:67:ARG:HB2	2.12	0.49
52:BV:22:VAL:HB	52:BV:94:LEU:CB	2.39	0.49
53:BW:29:LEU:HD23	53:BW:29:LEU:C	2.33	0.49
55:BY:87:LYS:C	55:BY:88:LYS:HD2	2.32	0.49
1:CA:163:C:O2'	1:CA:164:U:H5'	2.12	0.49
1:CA:447:G:N2	1:CA:488:C:N4	2.60	0.49
1:CA:498:U:C2'	1:CA:499:A:H5'	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:586:C:H1'	1:CA:878:G:O2'	2.12	0.49
2:CB:20:GLU:OE1	2:CB:20:GLU:HA	2.13	0.49
4:CD:149:ALA:O	4:CD:150:GLU:C	2.51	0.49
4:CD:154:ASN:HA	4:CD:159:ARG:HE	1.76	0.49
4:CD:158:ILE:O	4:CD:162:LEU:HG	2.13	0.49
4:CD:194:LEU:N	4:CD:194:LEU:CD2	2.75	0.49
4:CD:79:PHE:CD2	4:CD:79:PHE:C	2.86	0.49
5:CE:50:GLU:CD	5:CE:51:VAL:H	2.16	0.49
5:CE:60:TYR:HE2	5:CE:64:ARG:HE	1.60	0.49
8:CH:8:ASP:O	8:CH:12:ARG:HG3	2.12	0.49
10:CJ:78:ASN:C	10:CJ:80:LYS:N	2.63	0.49
12:CL:90:VAL:CG1	12:CL:93:LEU:HG	2.43	0.49
15:CO:37:ASN:N	15:CO:37:ASN:ND2	2.57	0.49
1:CA:472:A:H4'	16:CP:82:GLN:HE22	1.78	0.49
20:CT:14:LYS:HE3	20:CT:18:GLN:HE21	1.78	0.49
20:CT:53:LEU:O	20:CT:54:LYS:C	2.50	0.49
25:CY:16:LYS:O	25:CY:18:LEU:N	2.45	0.49
25:CY:9:GLU:O	25:CY:13:HIS:ND1	2.45	0.49
32:D6:28:ARG:HG2	32:D6:28:ARG:HH11	1.77	0.49
34:D8:30:ARG:HG2	35:DA:2393:A:H5''	1.95	0.49
35:DA:1141:U:C5'	35:DA:1142(A):A:O4'	2.60	0.49
35:DA:1166:C:H2'	35:DA:1167:U:C6	2.47	0.49
35:DA:1259:G:H2'	35:DA:1260:G:H8	1.76	0.49
35:DA:1532:C:O2	35:DA:1532:C:H2'	2.12	0.49
35:DA:1591:G:C2'	35:DA:1592:C:H5'	2.43	0.49
35:DA:1817:G:C2'	35:DA:1818:U:H5'	2.41	0.49
35:DA:184:C:H2'	35:DA:185:U:H6	1.76	0.49
35:DA:2052:G:N2	39:DE:149:ARG:HA	2.28	0.49
27:D1:37:ILE:HD12	35:DA:2080:G:OP1	2.12	0.49
35:DA:2124:G:C2'	35:DA:2125:G:H5'	2.43	0.49
35:DA:2292:C:H2'	35:DA:2293:C:C6	2.48	0.49
35:DA:2306:C:H5'	35:DA:2307:G:O5'	2.12	0.49
35:DA:2346:A:C2	35:DA:2383:G:N3	2.81	0.49
35:DA:2476:A:C2	35:DA:2477:C:C6	3.01	0.49
35:DA:2596:U:H2'	35:DA:2597:G:O4'	2.12	0.49
35:DA:2714:G:H5''	35:DA:2714:G:H8	1.77	0.49
35:DA:285:C:H2'	35:DA:286:C:C4'	2.42	0.49
35:DA:623:G:H2'	35:DA:624:C:C6	2.47	0.49
35:DA:811:U:O2	35:DA:1251:C:C6	2.65	0.49
37:DC:56:GLN:HB2	37:DC:168:THR:HA	1.95	0.49
38:DD:31:LYS:HZ2	38:DD:31:LYS:HA	1.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:96:HIS:HA	38:DD:102:LYS:HG2	1.93	0.49
39:DE:60:ASN:O	39:DE:61:ARG:C	2.51	0.49
39:DE:2:LYS:NZ	39:DE:95:ILE:O	2.31	0.49
40:DF:41:LEU:HD11	40:DF:184:TYR:CE1	2.42	0.49
41:DG:166:ASP:O	41:DG:170:ARG:N	2.44	0.49
43:DI:10:GLU:OE1	43:DI:11:ASN:N	2.46	0.49
44:DN:41:ASP:N	51:DU:64:ARG:NH1	2.53	0.49
44:DN:61:ARG:HH11	44:DN:61:ARG:HG3	1.76	0.49
45:DO:87:ILE:HG13	45:DO:91:LEU:CD1	2.42	0.49
46:DP:14:LYS:O	46:DP:15:ARG:HB2	2.13	0.49
46:DP:70:GLN:HG3	46:DP:71:VAL:N	2.28	0.49
46:DP:70:GLN:HG3	46:DP:71:VAL:HG12	1.94	0.49
46:DP:83:VAL:HG13	46:DP:83:VAL:O	2.13	0.49
47:DQ:71:ASP:N	47:DQ:94:VAL:O	2.38	0.49
35:DA:1286:A:OP1	48:DR:105:ARG:HD2	2.13	0.49
48:DR:2:ARG:NE	48:DR:5:LYS:NZ	2.60	0.49
48:DR:34:ILE:HG22	48:DR:35:THR:N	2.28	0.49
52:DV:72:VAL:O	52:DV:73:SER:CB	2.60	0.49
1:AA:1320:C:H5'	19:AS:70:LYS:CE	2.43	0.49
1:AA:1338:G:O2'	1:AA:1339:A:H5'	2.13	0.49
1:AA:1350:A:H2'	1:AA:1351:U:C6	2.48	0.49
1:AA:1396:A:O4'	1:AA:1398:A:HI'	2.13	0.49
1:AA:163:C:O2'	1:AA:164:U:H5'	2.13	0.49
1:AA:338:A:H2'	1:AA:339:C:C6	2.47	0.49
1:AA:939:G:H2'	1:AA:940:C:H6	1.77	0.49
2:AB:112:VAL:C	2:AB:115:LEU:HB3	2.29	0.49
2:AB:118:LEU:HD11	2:AB:141:GLU:OE1	2.13	0.49
2:AB:213:LEU:O	2:AB:217:ARG:HG2	2.12	0.49
5:AE:33:VAL:CG2	5:AE:43:LEU:HD13	2.35	0.49
7:AG:42:ILE:HG23	7:AG:117:ALA:CB	2.43	0.49
7:AG:15:ASP:HB2	7:AG:23:VAL:HB	1.95	0.49
7:AG:42:ILE:HG23	7:AG:117:ALA:CA	2.42	0.49
7:AG:79:ARG:HH11	7:AG:79:ARG:HG3	1.78	0.49
9:AI:87:GLN:C	9:AI:89:ASN:H	2.16	0.49
12:AL:41:ARG:HB3	12:AL:41:ARG:NH1	2.20	0.49
12:AL:55:VAL:O	12:AL:56:ALA:HB2	2.12	0.49
15:AO:53:HIS:O	15:AO:56:LEU:N	2.45	0.49
17:AQ:32:TYR:O	17:AQ:34:LYS:N	2.45	0.49
23:AW:49:C:C2'	23:AW:60:A:H4'	2.43	0.49
25:AY:29:ARG:HA	25:AY:32:ARG:NE	2.28	0.49
28:B2:24:LEU:HD11	28:B2:28:LYS:NZ	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:B5:40:LYS:NZ	31:B5:46:CYS:H	2.10	0.49
34:B8:34:TRP:HZ3	34:B8:41:ILE:HG23	1.78	0.49
35:BA:1024:G:H8	35:BA:1024:G:O5'	1.95	0.49
35:BA:1131:G:H1'	35:BA:1132:A:H8	1.77	0.49
35:BA:1255:U:H5''	35:BA:1256:G:H5''	1.92	0.49
35:BA:1465:G:C4'	35:BA:1528:A:H8	2.25	0.49
35:BA:1465:G:H5'	35:BA:1528:A:H8	1.78	0.49
35:BA:2336:A:H3'	35:BA:2337:G:H8	1.78	0.49
35:BA:2538:C:HO2'	35:BA:2539:C:H5'	1.76	0.49
35:BA:2677:G:C4	35:BA:2678:C:C5	3.00	0.49
35:BA:2684:U:H2'	35:BA:2685:G:O4'	2.12	0.49
35:BA:2864:G:H2'	35:BA:2865:U:O4'	2.13	0.49
35:BA:2892:A:N6	35:BA:2893:G:H21	2.11	0.49
35:BA:300:A:H5''	55:BY:97:ARG:NH1	2.27	0.49
33:B7:37:LYS:HG3	35:BA:458:G:C8	2.47	0.49
35:BA:589:C:H42	35:BA:668:G:H1	1.60	0.49
35:BA:612:C:H2'	35:BA:613:G:O4'	2.12	0.49
35:BA:71:A:H4'	35:BA:72:U:C5'	2.42	0.49
35:BA:942:G:H1'	35:BA:1189:A:H2	1.78	0.49
35:BA:990:A:C6	35:BA:1186:G:H1'	2.48	0.49
35:BA:1902:C:O2'	38:BD:244:ARG:HB2	2.12	0.49
38:BD:65:ILE:HD11	38:BD:67:PHE:CD1	2.47	0.49
35:BA:2574:G:O2'	39:BE:143:ASN:HB3	2.13	0.49
40:BF:117:ARG:NH2	40:BF:189:THR:O	2.46	0.49
40:BF:184:TYR:HD2	40:BF:185:ASP:N	2.11	0.49
40:BF:99:TYR:CG	40:BF:99:TYR:O	2.66	0.49
43:BI:120:ILE:O	43:BI:121:LYS:HB3	2.12	0.49
43:BI:38:LEU:N	43:BI:38:LEU:HD12	2.10	0.49
44:BN:13:TRP:H	44:BN:13:TRP:HD1	1.59	0.49
45:BO:13:ASN:HD21	45:BO:96:THR:H	1.59	0.49
35:BA:626:U:C2	46:BP:105:LEU:HG	2.47	0.49
46:BP:59:LEU:HA	46:BP:61:ARG:CD	2.43	0.49
47:BQ:20:ALA:CB	47:BQ:98:LYS:HB3	2.43	0.49
35:BA:1276:A:H1'	48:BR:16:HIS:HE1	1.78	0.49
50:BT:106:SER:O	50:BT:107:ASP:CB	2.61	0.49
39:BE:11:MET:N	50:BT:8:LYS:HE3	2.27	0.49
51:BU:108:GLU:HA	51:BU:111:GLU:HG2	1.95	0.49
52:BV:18:LEU:HA	52:BV:97:LYS:NZ	2.28	0.49
56:BZ:139:VAL:HG12	56:BZ:141:VAL:H	1.78	0.49
1:CA:1108:G:N3	1:CA:1108:G:H2'	2.27	0.49
1:CA:1401:G:H2'	1:CA:1402:C:C5'	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:570:G:C6	1:CA:873:A:N1	2.81	0.49
3:CC:150:LYS:HA	3:CC:169:ALA:HB2	1.93	0.49
3:CC:178:LEU:HD22	3:CC:178:LEU:H	1.78	0.49
4:CD:108:LEU:O	4:CD:110:PHE:HD1	1.95	0.49
5:CE:8:GLU:CA	5:CE:34:VAL:HG22	2.39	0.49
13:CM:54:VAL:O	13:CM:58:GLU:N	2.45	0.49
15:CO:85:LEU:HB2	15:CO:87:ILE:HD11	1.94	0.49
17:CQ:92:ARG:O	17:CQ:94:ASN:N	2.45	0.49
25:CY:2:THR:CG2	25:CY:4:LYS:HB3	2.42	0.49
26:D0:20:ARG:CD	26:D0:20:ARG:N	2.76	0.49
26:D0:29:GLN:O	26:D0:66:VAL:HA	2.13	0.49
27:D1:23:LYS:HZ3	27:D1:23:LYS:HA	1.76	0.49
31:D5:40:LYS:CE	31:D5:46:CYS:N	2.70	0.49
32:D6:15:GLU:CG	32:D6:15:GLU:O	2.58	0.49
34:D8:5:LYS:HD2	34:D8:5:LYS:N	2.27	0.49
35:DA:1591:G:O2'	35:DA:1592:C:H5'	2.13	0.49
35:DA:1790:C:H5''	35:DA:1791:A:OP1	2.13	0.49
35:DA:1814:G:H2'	35:DA:1815:A:C8	2.48	0.49
35:DA:2248:C:H3'	35:DA:2249:U:H6	1.78	0.49
35:DA:2569:G:C2	35:DA:2570:G:C8	3.01	0.49
35:DA:307:G:H21	35:DA:330:A:H62	1.61	0.49
35:DA:433:C:H2'	35:DA:434:U:C6	2.48	0.49
35:DA:602:G:N3	35:DA:602:G:H2'	2.27	0.49
35:DA:814:C:H2'	35:DA:815:C:H6	1.78	0.49
35:DA:838:C:N4	35:DA:940:G:H1	2.10	0.49
35:DA:998:C:H42	35:DA:1157:G:H1	1.60	0.49
37:DC:99:ILE:HG23	37:DC:103:ILE:CB	2.43	0.49
38:DD:25:THR:HG21	38:DD:82:ILE:N	2.27	0.49
39:DE:116:VAL:HG22	39:DE:117:MET:H	1.78	0.49
39:DE:197:ILE:HG13	39:DE:199:ARG:NH1	2.24	0.49
41:DG:48:GLU:O	41:DG:49:ASP:CB	2.61	0.49
42:DH:40:GLU:O	42:DH:41:MET:HB2	2.12	0.49
42:DH:17:VAL:HB	42:DH:45:VAL:HG22	1.94	0.49
42:DH:67:LEU:C	42:DH:71:LEU:HD22	2.32	0.49
44:DN:71:ILE:HD13	44:DN:86:PRO:HA	1.94	0.49
44:DN:93:THR:C	44:DN:94:HIS:CG	2.85	0.49
51:DU:98:LEU:C	51:DU:100:VAL:N	2.65	0.49
56:DZ:14:LYS:C	56:DZ:16:SER:N	2.66	0.49
56:DZ:58:VAL:O	56:DZ:58:VAL:HG12	2.13	0.49
1:AA:1069:C:O2'	1:AA:1192:C:H1'	2.12	0.49
1:AA:1459:C:H2'	1:AA:1460:A:O4'	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:596:C:H2'	1:AA:596:C:O2	2.13	0.49
1:AA:648:A:H2'	1:AA:649:G:C8	2.47	0.49
1:AA:574:A:N3	1:AA:883:C:H1'	2.28	0.49
1:AA:972:C:C5'	10:AJ:57:LYS:HG3	2.43	0.49
2:AB:100:GLY:O	2:AB:104:ASN:N	2.46	0.49
3:AC:60:ALA:CB	3:AC:63:ASN:HD21	2.25	0.49
4:AD:52:SER:OG	4:AD:55:ALA:N	2.38	0.49
5:AE:82:VAL:CG2	5:AE:138:ALA:HA	2.39	0.49
5:AE:72:GLN:HE22	5:AE:77:PRO:CD	2.24	0.49
5:AE:80:ILE:CD1	5:AE:91:LEU:HB2	2.43	0.49
5:AE:89:ILE:HG12	5:AE:90:VAL:N	2.26	0.49
8:AH:47:GLY:O	8:AH:62:TYR:N	2.36	0.49
11:AK:58:PRO:HA	11:AK:90:GLY:HA3	1.95	0.49
12:AL:21:LYS:HD2	12:AL:21:LYS:N	2.21	0.49
17:AQ:82:MET:HA	17:AQ:85:VAL:HG23	1.95	0.49
17:AQ:86:GLU:HA	17:AQ:89:LEU:HB3	1.94	0.49
18:AR:34:TYR:O	18:AR:35:ARG:HG2	2.12	0.49
20:AT:73:HIS:C	20:AT:74:LYS:HD3	2.33	0.49
20:AT:97:ALA:O	20:AT:99:LEU:N	2.42	0.49
25:AY:126:ARG:HA	25:AY:169:ILE:CD1	2.43	0.49
25:AY:29:ARG:NH2	25:AY:88:LEU:HA	2.28	0.49
25:AY:70:SER:CB	25:AY:76:LEU:HB2	2.42	0.49
25:AY:73:GLN:HG3	25:AY:74:ASN:OD1	2.12	0.49
26:B0:36:ILE:HG13	26:B0:36:ILE:O	2.11	0.49
29:B3:26:LEU:O	29:B3:27:GLY:C	2.50	0.49
35:BA:1277:G:H2'	35:BA:1278:A:C8	2.48	0.49
35:BA:1360:A:C6	35:BA:1372:U:C5	3.00	0.49
35:BA:769:G:H4'	35:BA:1379:A:N1	2.27	0.49
35:BA:1493:C:C2'	35:BA:1493:C:O2	2.61	0.49
35:BA:2170:A:H5'	35:BA:2171:A:OP2	2.13	0.49
35:BA:2543:G:H2'	35:BA:2544:G:H8	1.75	0.49
35:BA:2880:C:O3'	48:BR:90:ARG:NH1	2.44	0.49
35:BA:694:U:C2'	35:BA:695:G:O5'	2.61	0.49
35:BA:862:G:H3'	35:BA:863:A:H8	1.78	0.49
35:BA:917:A:H2'	35:BA:918:A:O4'	2.12	0.49
37:BC:44:HIS:HD2	37:BC:175:VAL:CA	2.26	0.49
35:BA:2591:C:P	38:BD:239:ARG:HG3	2.53	0.49
38:BD:27:THR:HG21	38:BD:83:GLU:CG	2.28	0.49
35:BA:1659:U:OP2	39:BE:132:HIS:CE1	2.66	0.49
40:BF:121:GLY:O	40:BF:123:LEU:N	2.46	0.49
42:BH:113:VAL:HG21	42:BH:151:ILE:CG2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BI:36:ALA:O	43:BI:37:VAL:HG23	2.13	0.49
43:BI:82:ARG:HG3	43:BI:82:ARG:NH1	2.27	0.49
45:BO:1:MET:H3	45:BO:1:MET:HE2	1.78	0.49
46:BP:138:LEU:O	46:BP:138:LEU:HD22	2.12	0.49
48:BR:101:ALA:HB1	53:BW:38:TYR:HE1	1.78	0.49
48:BR:56:LYS:HE2	48:BR:94:TYR:CE2	2.48	0.49
49:BS:46:VAL:HG12	49:BS:47:THR:N	2.27	0.49
49:BS:28:VAL:N	49:BS:89:ARG:HG2	2.28	0.49
52:BV:19:LYS:NZ	52:BV:20:LEU:N	2.56	0.49
54:BX:14:SER:O	54:BX:17:ALA:N	2.45	0.49
54:BX:26:TYR:H	54:BX:26:TYR:HD1	1.61	0.49
54:BX:62:LYS:CB	54:BX:69:TYR:H	2.26	0.49
56:BZ:58:VAL:CA	56:BZ:68:PRO:HA	2.42	0.49
56:BZ:60:GLU:HA	56:BZ:66:SER:CB	2.41	0.49
1:CA:1350:A:H2'	1:CA:1351:U:C6	2.48	0.49
1:CA:16:A:N1	1:CA:919:A:C2	2.81	0.49
1:CA:189(B):C:H2'	1:CA:189(C):C:C6	2.48	0.49
1:CA:199:G:H2'	1:CA:200:G:H8	1.77	0.49
1:CA:334:C:C2'	1:CA:335:C:H5'	2.42	0.49
1:CA:376:G:O2'	1:CA:377:G:H5'	2.13	0.49
1:CA:659:U:H2'	1:CA:660:G:C8	2.48	0.49
1:CA:665:A:C2'	1:CA:725:G:N2	2.72	0.49
1:CA:777:A:H2'	1:CA:778:G:H8	1.77	0.49
1:CA:838:G:O2'	1:CA:839:U:H5''	2.13	0.49
1:CA:863:U:H2'	1:CA:865:A:OP2	2.13	0.49
2:CB:115:LEU:HA	2:CB:145:LEU:HD11	1.95	0.49
2:CB:97:TRP:HH2	2:CB:176:GLU:CB	2.25	0.49
4:CD:150:GLU:HG2	4:CD:151:LYS:H	1.77	0.49
4:CD:65:ARG:HB2	4:CD:75:PHE:CZ	2.48	0.49
7:CG:50:ILE:HG21	7:CG:61:VAL:HG21	1.94	0.49
8:CH:114:THR:HG21	8:CH:119:LEU:HD21	1.95	0.49
8:CH:27:PRO:O	8:CH:32:LYS:NZ	2.41	0.49
13:CM:112:GLY:C	13:CM:113:PRO:HG2	2.33	0.49
17:CQ:86:GLU:HB3	17:CQ:90:ILE:HG12	1.93	0.49
20:CT:73:HIS:C	20:CT:74:LYS:HD3	2.32	0.49
21:CU:18:TYR:HA	21:CU:22:ARG:HD3	1.95	0.49
23:CW:43:G:C3'	23:CW:44:A:H5''	2.43	0.49
27:D1:72:GLU:O	27:D1:76:ARG:NH1	2.46	0.49
27:D1:89:GLU:O	27:D1:93:GLU:CD	2.50	0.49
27:D1:89:GLU:HG2	27:D1:90:ILE:CD1	2.42	0.49
28:D2:12:GLU:O	28:D2:14:ARG:CD	2.61	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D2:32:LEU:HG	28:D2:33:MET:SD	2.52	0.49
29:D3:52:HIS:CD2	36:DB:83:G:C4'	2.95	0.49
35:DA:1516:C:H2'	35:DA:1517:G:H8	1.77	0.49
35:DA:1515:G:H2'	35:DA:1516:C:H5'	1.93	0.49
35:DA:2030:A:H5''	35:DA:2031:A:OP1	2.13	0.49
35:DA:2250:G:H8	35:DA:2496:C:H5''	1.77	0.49
35:DA:2442:C:O2'	35:DA:2443:C:H5'	2.12	0.49
35:DA:2640:G:H5'	35:DA:2640:G:H8	1.78	0.49
35:DA:2676:C:O2'	35:DA:2677:G:H5'	2.13	0.49
35:DA:375:C:H2'	35:DA:376:C:C6	2.48	0.49
35:DA:32:C:H42	35:DA:473:G:H1	1.60	0.49
35:DA:613:G:C6	35:DA:615:G:O6	2.66	0.49
35:DA:61:G:C2'	35:DA:62:C:H5'	2.43	0.49
35:DA:66:C:O2'	35:DA:67:U:H5'	2.12	0.49
35:DA:773:U:H4'	38:DD:47:GLY:HA2	1.94	0.49
35:DA:869:G:H1'	47:DQ:8:LYS:HZ3	1.75	0.49
37:DC:64:LEU:HD12	37:DC:66:HIS:HB2	1.95	0.49
38:DD:117:VAL:CG2	38:DD:118:VAL:H	2.24	0.49
38:DD:264:LYS:CE	38:DD:266:SER:HB2	2.43	0.49
40:DF:39:TRP:CB	40:DF:101:LEU:HD22	2.42	0.49
42:DH:103:LEU:HG	42:DH:105:LEU:CD1	2.43	0.49
42:DH:18:GLU:CB	42:DH:25:LYS:HD2	2.43	0.49
43:DI:54:GLN:OE1	43:DI:54:GLN:C	2.51	0.49
44:DN:31:ALA:O	44:DN:32:THR:C	2.51	0.49
46:DP:39:LYS:C	46:DP:41:ARG:N	2.66	0.49
47:DQ:141:GLN:HE21	56:DZ:72:ARG:CA	2.23	0.49
48:DR:28:LEU:C	48:DR:28:LEU:HD13	2.32	0.49
48:DR:60:LEU:O	48:DR:63:ARG:N	2.46	0.49
48:DR:6:SER:HA	48:DR:8:ARG:HH21	1.78	0.49
49:DS:83:LYS:HA	49:DS:104:GLY:CA	2.40	0.49
51:DU:79:PHE:O	51:DU:83:LEU:HD13	2.12	0.49
53:DW:5:ALA:HB1	53:DW:50:VAL:HG22	1.95	0.49
54:DX:46:ALA:C	54:DX:47:PHE:CD1	2.86	0.49
1:AA:1128:C:H5'	9:AI:16:ARG:NH1	2.28	0.49
1:AA:1163:C:H2'	1:AA:1164:G:C8	2.46	0.49
1:AA:1168:A:C2	1:AA:1169:A:C4	3.01	0.49
1:AA:961:U:OP2	1:AA:1223:C:H4'	2.13	0.49
1:AA:1416:G:C4	1:AA:1417:G:C8	3.00	0.49
1:AA:243:A:O2'	1:AA:244:U:OP2	2.30	0.49
1:AA:309:G:H2'	1:AA:310:G:H8	1.78	0.49
1:AA:577:G:C4	1:AA:816:A:C2	3.00	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:645:C:O2'	1:AA:646:U:H5'	2.12	0.49
1:AA:858:G:O6	1:AA:869:G:C8	2.66	0.49
1:AA:972:C:OP2	10:AJ:57:LYS:HE2	2.13	0.49
2:AB:9:GLU:O	2:AB:13:ALA:CB	2.60	0.49
2:AB:69:LEU:HD13	2:AB:91:PRO:O	2.13	0.49
4:AD:108:LEU:O	4:AD:110:PHE:HD1	1.96	0.49
4:AD:110:PHE:H	4:AD:110:PHE:HD1	1.59	0.49
6:AF:10:LEU:HD21	6:AF:22:GLU:HB3	1.94	0.49
7:AG:40:ALA:O	7:AG:44:TYR:HD1	1.96	0.49
8:AH:27:PRO:O	8:AH:32:LYS:NZ	2.44	0.49
8:AH:6:ILE:N	8:AH:6:ILE:HD12	2.27	0.49
11:AK:33:THR:C	11:AK:40:ILE:HG12	2.33	0.49
11:AK:15:ALA:HA	11:AK:76:GLY:O	2.13	0.49
1:AA:1227:A:O2'	13:AM:115:LYS:HB2	2.13	0.49
16:AP:60:LEU:O	16:AP:62:VAL:N	2.46	0.49
23:AW:41:C:H2'	23:AW:42:C:H6	1.77	0.49
25:AY:34:ASN:OD1	25:AY:35:PRO:HD2	2.12	0.49
25:AY:76:LEU:HD11	25:AY:98:ALA:C	2.33	0.49
27:B1:20:ARG:NH1	27:B1:41:ARG:CZ	2.72	0.49
28:B2:53:LEU:HA	28:B2:56:GLN:HG3	1.95	0.49
28:B2:14:ARG:HD2	28:B2:57:ILE:HD12	1.95	0.49
35:BA:1024:G:N2	35:BA:1142(A):A:H2	2.11	0.49
35:BA:1567:A:H3'	38:BD:86:PRO:HG3	1.94	0.49
35:BA:1682:G:H2'	35:BA:1683:C:C6	2.48	0.49
35:BA:49:A:C5	35:BA:177:G:C6	3.01	0.49
35:BA:1926:U:H2'	35:BA:1928:A:OP2	2.13	0.49
35:BA:2225:A:H1'	35:BA:2226:C:OP2	2.13	0.49
35:BA:2364:C:C2'	35:BA:2365:G:H5'	2.43	0.49
35:BA:2497:A:C8	35:BA:2497:A:O5'	2.66	0.49
35:BA:256:A:H2'	35:BA:257:A:H8	1.77	0.49
35:BA:2708:G:H1'	48:BR:71:GLN:OE1	2.13	0.49
37:BC:41:VAL:HB	37:BC:178:ALA:CB	2.42	0.49
38:BD:264:LYS:CE	38:BD:266:SER:HB2	2.43	0.49
39:BE:60:ASN:O	39:BE:63:LEU:N	2.45	0.49
40:BF:32:LEU:O	40:BF:35:GLU:N	2.46	0.49
41:BG:7:LEU:HG	41:BG:104:GLU:HG2	1.93	0.49
41:BG:73:ALA:CB	41:BG:87:PRO:HG2	2.38	0.49
27:B1:71:TYR:CE1	43:BI:27:ARG:HG3	2.48	0.49
43:BI:54:GLN:HA	43:BI:57:ARG:HH11	1.77	0.49
45:BO:71:ARG:NH1	45:BO:71:ARG:HG3	2.28	0.49
45:BO:77:ILE:HD11	50:BT:72:VAL:HG13	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:23:PRO:HB2	46:BP:33:ARG:NE	2.27	0.49
35:BA:2406:U:N3	46:BP:72:PRO:HD2	2.27	0.49
46:BP:97:PRO:C	46:BP:99:LEU:N	2.67	0.49
47:BQ:25:ASP:CB	47:BQ:67:ARG:HH22	2.26	0.49
49:BS:66:ALA:HB1	49:BS:98:VAL:O	2.13	0.49
49:BS:95:HIS:O	49:BS:97:ARG:N	2.46	0.49
51:BU:18:LEU:HD21	51:BU:22:LYS:HE2	1.95	0.49
51:BU:98:LEU:HD22	52:BV:2:PHE:HZ	1.78	0.49
52:BV:72:VAL:HG13	52:BV:88:ARG:NH2	2.28	0.49
52:BV:69:LYS:HB2	52:BV:93:GLU:OE2	2.13	0.49
53:BW:5:ALA:HB1	53:BW:50:VAL:HG22	1.94	0.49
53:BW:92:ARG:HH11	53:BW:92:ARG:HG2	1.78	0.49
54:BX:83:VAL:C	54:BX:85:PRO:HD3	2.33	0.49
55:BY:38:ILE:CG2	55:BY:39:VAL:N	2.76	0.49
55:BY:14:LEU:O	55:BY:72:VAL:HA	2.13	0.49
56:BZ:121:HIS:C	56:BZ:123:ASP:H	2.16	0.49
56:BZ:129:SER:C	56:BZ:131:ARG:H	2.15	0.49
56:BZ:73:GLN:O	56:BZ:87:ASP:OD1	2.31	0.49
1:CA:1104:G:H2'	1:CA:1105:A:C8	2.48	0.49
1:CA:498:U:H2'	1:CA:499:A:C5'	2.42	0.49
1:CA:894:G:H2'	1:CA:895:G:H8	1.78	0.49
1:CA:908:A:C2	1:CA:909:A:N7	2.81	0.49
1:CA:90:U:O5'	1:CA:90:U:H6	1.96	0.49
3:CC:91:LEU:CB	3:CC:99:VAL:HG21	2.43	0.49
4:CD:18:LYS:NZ	4:CD:31:CYS:CB	2.76	0.49
4:CD:96:LEU:HD13	4:CD:96:LEU:N	2.28	0.49
5:CE:130:ASN:O	5:CE:131:ILE:C	2.49	0.49
7:CG:112:PRO:HD2	7:CG:113:GLU:OE2	2.12	0.49
7:CG:11:GLN:HG3	7:CG:12:LEU:N	2.28	0.49
8:CH:60:ARG:HH11	8:CH:60:ARG:HG3	1.78	0.49
10:CJ:50:ILE:HD11	14:CN:41:ARG:CD	2.38	0.49
10:CJ:55:LYS:O	10:CJ:56:HIS:CG	2.66	0.49
12:CL:7:ILE:O	12:CL:10:LEU:HB2	2.13	0.49
12:CL:84:LEU:HD22	12:CL:101:VAL:HG21	1.95	0.49
13:CM:86:CYS:O	13:CM:89:GLY:N	2.45	0.49
1:CA:1308:U:OP1	13:CM:98:VAL:N	2.46	0.49
15:CO:66:LEU:O	15:CO:69:TYR:HB3	2.12	0.49
15:CO:70:LEU:CD2	15:CO:78:TYR:HA	2.42	0.49
18:CR:87:ARG:HH11	18:CR:87:ARG:HB3	1.77	0.49
19:CS:16:LEU:N	19:CS:16:LEU:CD1	2.75	0.49
25:CY:173:ASP:C	25:CY:175:LEU:N	2.67	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:D1:19:GLN:OE1	27:D1:44:PRO:HB3	2.13	0.49
28:D2:14:ARG:C	28:D2:14:ARG:NE	2.59	0.49
35:DA:1221:C:H2'	35:DA:1221(A):C:C6	2.42	0.49
35:DA:1250:G:H3'	35:DA:1251:C:H5'	1.94	0.49
35:DA:1278:A:OP1	48:DR:36:THR:HG22	2.13	0.49
35:DA:1310:G:H1	35:DA:1604:C:H42	1.60	0.49
35:DA:1681:G:O2'	35:DA:1762:A:H2'	2.13	0.49
35:DA:1800:C:C4	35:DA:1818:U:O2	2.66	0.49
35:DA:1887:C:C3'	35:DA:1888:G:H5''	2.42	0.49
35:DA:2247:A:O2'	35:DA:2248:C:H5'	2.13	0.49
35:DA:2251:G:N3	35:DA:2450:A:H1'	2.28	0.49
35:DA:2843:G:C2	35:DA:2875:C:N3	2.81	0.49
35:DA:2854:G:H1	35:DA:2863:C:H42	1.61	0.49
27:D1:34:THR:CG2	35:DA:387:U:O3'	2.61	0.49
35:DA:609:A:H2'	35:DA:610:G:O4'	2.13	0.49
35:DA:612:C:H2'	35:DA:613:G:O4'	2.13	0.49
35:DA:707:G:H3'	35:DA:708:C:H6	1.78	0.49
35:DA:785:G:H2'	35:DA:786:C:H6	1.76	0.49
36:DB:29:A:H2'	36:DB:30:C:C6	2.48	0.49
36:DB:52:A:O2'	36:DB:53:A:H8	1.96	0.49
37:DC:50:ASP:O	37:DC:52:ARG:N	2.45	0.49
37:DC:76:ALA:C	37:DC:78:ALA:H	2.16	0.49
39:DE:181:LEU:O	39:DE:182:LEU:CB	2.60	0.49
39:DE:78:LEU:C	39:DE:79:ARG:HD2	2.34	0.49
41:DG:114:ILE:HD12	41:DG:117:PHE:CG	2.48	0.49
41:DG:131:TYR:CD1	41:DG:132:ASN:N	2.81	0.49
42:DH:66:GLY:CA	42:DH:69:ARG:HB2	2.35	0.49
44:DN:52:VAL:O	44:DN:120:LEU:HD22	2.12	0.49
35:DA:1141:U:OP2	44:DN:63:THR:HG21	2.13	0.49
46:DP:6:LEU:O	46:DP:6:LEU:HD23	2.12	0.49
47:DQ:20:ALA:HA	47:DQ:98:LYS:HD3	1.94	0.49
48:DR:103:ARG:HB3	48:DR:109:ALA:O	2.13	0.49
52:DV:15:GLU:CB	52:DV:16:PRO:HD2	2.41	0.49
56:DZ:48:PHE:HE2	56:DZ:71:VAL:HG11	1.73	0.49
1:AA:1308:U:OP1	13:AM:98:VAL:N	2.45	0.48
1:AA:1435:G:C2	1:AA:1436:U:N3	2.81	0.48
1:AA:390:C:H2'	1:AA:391:G:H8	1.78	0.48
1:AA:729:A:H2'	1:AA:730:G:H8	1.78	0.48
4:AD:173:TRP:O	4:AD:186:LEU:HB2	2.12	0.48
5:AE:60:TYR:HE2	5:AE:64:ARG:HE	1.60	0.48
6:AF:85:VAL:HG12	6:AF:85:VAL:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:50:ARG:HG2	8:AH:50:ARG:HH11	1.77	0.48
9:AI:28:VAL:CG1	9:AI:29:ASN:N	2.75	0.48
7:AG:16:LEU:CD1	9:AI:41:VAL:HG12	2.43	0.48
9:AI:53:VAL:HG12	9:AI:95:LYS:HE3	1.95	0.48
12:AL:32:PHE:CB	12:AL:84:LEU:HD21	2.43	0.48
13:AM:48:LEU:HD11	13:AM:53:VAL:HG22	1.95	0.48
15:AO:11:VAL:HG21	15:AO:34:LEU:CD2	2.43	0.48
15:AO:78:TYR:O	15:AO:82:ILE:HG22	2.13	0.48
16:AP:60:LEU:HD23	16:AP:64:ALA:CB	2.30	0.48
1:AA:235:C:H5'	17:AQ:70:ARG:HG2	1.94	0.48
19:AS:9:VAL:O	19:AS:10:PHE:C	2.51	0.48
20:AT:29:LYS:O	20:AT:32:ALA:HB3	2.12	0.48
23:AW:26:C:H2'	23:AW:27:G:O4'	2.12	0.48
23:AW:38:A:H2'	23:AW:39:A:H5''	1.95	0.48
26:B0:32:ARG:N	26:B0:35:ASN:ND2	2.47	0.48
27:B1:90:ILE:HD13	27:B1:90:ILE:N	2.27	0.48
35:BA:1788:C:O5'	35:BA:1788:C:H6	1.95	0.48
35:BA:189:G:C4	35:BA:205:G:N2	2.80	0.48
35:BA:1916:A:H3'	35:BA:1917:U:H6	1.78	0.48
35:BA:1937:A:H2'	35:BA:1938:A:H5'	1.94	0.48
35:BA:2115:G:H22	35:BA:2170:A:N6	2.11	0.48
35:BA:2303:G:H2'	35:BA:2304:G:O4'	2.13	0.48
35:BA:2305:A:N6	41:BG:43:LEU:HA	2.28	0.48
35:BA:2547:U:H2'	35:BA:2548:G:H8	1.78	0.48
35:BA:1953:A:H2	35:BA:2549:G:N3	2.11	0.48
35:BA:2562:U:H2'	35:BA:2563:U:C5'	2.43	0.48
35:BA:2673:G:HO2'	35:BA:2674:G:H5'	1.77	0.48
35:BA:285:C:H2'	35:BA:286:C:C4'	2.43	0.48
35:BA:66:C:O2'	35:BA:67:U:H5'	2.13	0.48
36:BB:83:G:C2	36:BB:84:C:C6	3.01	0.48
38:BD:201:HIS:C	38:BD:203:ASN:N	2.66	0.48
38:BD:267:SER:O	38:BD:269:PHE:HD1	1.95	0.48
38:BD:34:VAL:O	38:BD:34:VAL:CG1	2.61	0.48
39:BE:55:ASN:ND2	39:BE:75:VAL:HG22	2.28	0.48
40:BF:6:VAL:CG2	40:BF:124:LEU:HA	2.41	0.48
40:BF:1:MET:O	40:BF:2:LYS:C	2.50	0.48
41:BG:121:ASN:HB2	41:BG:181:ARG:NH2	2.28	0.48
41:BG:9:ARG:HG3	41:BG:9:ARG:HH11	1.77	0.48
43:BI:29:TYR:CD2	43:BI:30:LEU:HD23	2.47	0.48
44:BN:125:GLY:HA3	44:BN:126:PRO:O	2.12	0.48
44:BN:56:ASN:HA	44:BN:124:ALA:CA	2.40	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:4:PRO:HA	45:BO:21:CYS:SG	2.53	0.48
46:BP:23:PRO:HB2	46:BP:33:ARG:CG	2.42	0.48
46:BP:48:PRO:HG2	46:BP:49:ARG:N	2.28	0.48
47:BQ:38:GLU:HB2	47:BQ:127:ILE:CG2	2.43	0.48
47:BQ:88:GLY:O	47:BQ:89:ASN:CB	2.61	0.48
48:BR:16:HIS:O	48:BR:19:ALA:N	2.44	0.48
48:BR:3:HIS:HB2	48:BR:4:LEU:H	1.47	0.48
50:BT:100:TYR:CD2	50:BT:103:ARG:CZ	2.93	0.48
50:BT:52:ILE:HG22	50:BT:61:PHE:HB3	1.95	0.48
45:BO:76:ALA:CB	50:BT:75:ILE:HD13	2.42	0.48
51:BU:33:ARG:O	51:BU:36:ARG:N	2.46	0.48
52:BV:72:VAL:HA	52:BV:88:ARG:HH22	1.78	0.48
55:BY:14:LEU:HD12	55:BY:15:VAL:N	2.24	0.48
55:BY:7:VAL:HG23	55:BY:8:LYS:HD2	1.94	0.48
56:BZ:139:VAL:C	56:BZ:141:VAL:N	2.67	0.48
56:BZ:16:SER:O	56:BZ:19:ARG:HG2	2.13	0.48
1:CA:1036:G:H2'	1:CA:1036:G:N3	2.27	0.48
1:CA:1049:U:H1'	1:CA:1201:A:C8	2.48	0.48
1:CA:1349:A:O2'	1:CA:1350:A:H5'	2.13	0.48
1:CA:1516:G:N2	1:CA:1518:A:H3'	2.28	0.48
1:CA:324:G:N2	1:CA:327:A:C8	2.81	0.48
1:CA:645:C:H2'	1:CA:646:U:C6	2.47	0.48
1:CA:895:G:C6	1:CA:896:C:N4	2.81	0.48
4:CD:2:GLY:O	4:CD:3:ARG:C	2.51	0.48
4:CD:34:GLU:O	4:CD:35:ARG:HG2	2.13	0.48
4:CD:59:ARG:NH2	4:CD:66:ARG:HH22	2.05	0.48
5:CE:91:LEU:HD23	5:CE:110:LEU:HD11	1.95	0.48
8:CH:114:THR:HG22	8:CH:117:GLY:O	2.13	0.48
8:CH:68:ARG:HG3	8:CH:69:ARG:N	2.23	0.48
9:CI:17:VAL:CG1	9:CI:81:ILE:HD13	2.43	0.48
11:CK:66:LEU:O	11:CK:67:ASP:C	2.49	0.48
12:CL:32:PHE:HD1	12:CL:86:ARG:HA	1.78	0.48
13:CM:10:PRO:CB	13:CM:18:ALA:HB1	2.42	0.48
16:CP:80:PHE:O	16:CP:81:ARG:C	2.51	0.48
17:CQ:45:HIS:O	17:CQ:73:VAL:HG23	2.13	0.48
23:CW:11:A:H2'	23:CW:12:G:C8	2.48	0.48
25:CY:108:GLU:HG2	25:CY:111:ARG:HH21	1.78	0.48
25:CY:123:GLU:HA	25:CY:126:ARG:HH11	1.78	0.48
25:CY:26:ALA:HA	25:CY:179:LYS:NZ	2.28	0.48
25:CY:93:SER:C	25:CY:95:LYS:H	2.16	0.48
27:D1:13:ILE:CG1	27:D1:14:VAL:H	2.04	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:D1:76:ARG:CA	27:D1:78:LYS:HZ3	2.25	0.48
29:D3:13:ILE:HD12	35:DA:989:G:N7	2.28	0.48
29:D3:26:LEU:HD21	29:D3:46:ASN:HB3	1.95	0.48
35:DA:1021:A:C3'	35:DA:1021:A:C8	2.96	0.48
35:DA:1213:A:N3	35:DA:1238:G:H1'	2.28	0.48
35:DA:1445(A):C:H2'	35:DA:1446:C:C6	2.48	0.48
35:DA:1553:A:N7	35:DA:1555:G:C5	2.81	0.48
35:DA:1637:A:H2'	35:DA:1638:C:C6	2.47	0.48
35:DA:154(A):C:H5	35:DA:171:G:H1	1.57	0.48
35:DA:1848:A:H2'	35:DA:1849:G:C8	2.47	0.48
35:DA:199:A:N3	35:DA:2433:A:C2	2.81	0.48
35:DA:2019:A:N1	35:DA:2020:A:C4	2.81	0.48
35:DA:204:A:H5'	35:DA:206:U:O4'	2.13	0.48
35:DA:2251:G:H2'	35:DA:2252:G:O4'	2.13	0.48
35:DA:2350:C:H2'	35:DA:2351:G:O4'	2.13	0.48
35:DA:2517:C:H2'	35:DA:2542:A:H2	1.78	0.48
35:DA:2566:A:N6	45:DO:28:SER:HB2	2.28	0.48
35:DA:2619:C:O2'	35:DA:2620:C:H5'	2.13	0.48
35:DA:286:C:O2	35:DA:286:C:C2'	2.58	0.48
35:DA:239:U:O2'	35:DA:622:G:H4'	2.13	0.48
28:D2:51:ARG:CZ	35:DA:72:U:H5'	2.43	0.48
35:DA:847:U:C2'	35:DA:848:G:H5''	2.30	0.48
36:DB:18:G:H2'	36:DB:19:G:H8	1.78	0.48
38:DD:133:LEU:HB2	38:DD:187:GLY:HA2	1.95	0.48
38:DD:201:HIS:C	38:DD:203:ASN:N	2.65	0.48
35:DA:1567:A:N6	38:DD:21:PHE:CE2	2.81	0.48
38:DD:239:ARG:NH2	38:DD:239:ARG:HG3	2.26	0.48
39:DE:47:VAL:HG12	39:DE:49:LEU:CD2	2.43	0.48
39:DE:48:GLN:O	39:DE:49:LEU:HD13	2.13	0.48
40:DF:103:LYS:O	40:DF:104:LYS:C	2.50	0.48
42:DH:85:LYS:HD3	42:DH:133:VAL:HB	1.94	0.48
42:DH:138:LYS:HA	42:DH:141:VAL:HB	1.95	0.48
43:DI:54:GLN:HA	43:DI:57:ARG:HH11	1.77	0.48
43:DI:91:SER:H	43:DI:121:LYS:HE2	1.77	0.48
43:DI:94:ALA:CB	43:DI:114:LEU:HD12	2.43	0.48
47:DQ:32:TYR:N	47:DQ:32:TYR:CD1	2.81	0.48
48:DR:14:SER:O	48:DR:15:SER:C	2.52	0.48
48:DR:39:PRO:O	48:DR:40:LYS:C	2.51	0.48
35:DA:1453:U:OP1	48:DR:77:ARG:NH1	2.46	0.48
49:DS:87:PHE:HB2	49:DS:106:ARG:HD3	1.94	0.48
49:DS:13:ARG:O	49:DS:15:ARG:HD3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:92:TYR:CD2	49:DS:97:ARG:NH1	2.80	0.48
50:DT:27:THR:HG23	50:DT:28:VAL:N	2.28	0.48
51:DU:74:LEU:CD1	51:DU:79:PHE:HB2	2.42	0.48
52:DV:4:ILE:HB	52:DV:40:LEU:HD11	1.95	0.48
52:DV:75:PHE:CE1	52:DV:89:GLN:HB2	2.48	0.48
52:DV:18:LEU:HA	52:DV:97:LYS:NZ	2.28	0.48
53:DW:17:VAL:O	53:DW:18:ARG:C	2.52	0.48
53:DW:86:LEU:HD12	53:DW:87:PRO:CD	2.43	0.48
56:DZ:102:LEU:HD12	56:DZ:121:HIS:O	2.12	0.48
1:AA:1447:A:OP2	1:AA:1452:C:C6	2.65	0.48
1:AA:321:A:N7	1:AA:328:C:O2	2.46	0.48
1:AA:372:C:H4'	1:AA:373:A:OP1	2.14	0.48
1:AA:738:C:H5''	6:AF:69:GLU:CB	2.42	0.48
1:AA:788:U:C5	1:AA:789:U:C5	3.01	0.48
1:AA:997:U:H2'	1:AA:998:G:C8	2.48	0.48
2:AB:19:HIS:CE1	2:AB:206:ASP:HB2	2.47	0.48
2:AB:80:ILE:CD1	2:AB:215:LEU:HD12	2.43	0.48
4:AD:159:ARG:HH11	4:AD:159:ARG:HG3	1.78	0.48
7:AG:85:TYR:CE1	7:AG:154:TYR:HE1	2.31	0.48
8:AH:28:ALA:HB2	8:AH:57:PRO:O	2.13	0.48
9:AI:126:SER:O	9:AI:127:LYS:CB	2.61	0.48
1:AA:1250:A:H4'	9:AI:68:GLY:O	2.13	0.48
12:AL:25:PRO:C	12:AL:27:LEU:H	2.15	0.48
14:AN:24:CYS:HB3	14:AN:27:CYS:O	2.12	0.48
19:AS:62:ILE:HD12	19:AS:63:THR:N	2.26	0.48
25:AY:61:PRO:HD3	25:AY:66:LEU:HA	1.95	0.48
32:B6:40:CYS:HB2	32:B6:46:HIS:CE1	2.45	0.48
35:BA:103:A:H2'	35:BA:104:U:C6	2.47	0.48
35:BA:1274:A:N3	35:BA:1297:C:H1'	2.27	0.48
35:BA:1348:G:H2'	35:BA:1349:A:C5'	2.30	0.48
35:BA:1354:A:C8	35:BA:1355:G:C8	3.01	0.48
35:BA:1399:C:H2'	35:BA:1400:G:C8	2.48	0.48
35:BA:1400:G:H2'	35:BA:1401:G:C8	2.47	0.48
35:BA:1446:C:O2'	35:BA:1447:G:H5'	2.12	0.48
35:BA:1982:C:H2'	35:BA:1983:C:H6	1.79	0.48
35:BA:2534:A:H2'	35:BA:2535:G:O4'	2.13	0.48
35:BA:2712:U:C1'	35:BA:2712(A):A:C8	2.95	0.48
35:BA:2736:G:O2'	35:BA:2737:G:H5'	2.13	0.48
35:BA:449:A:C2'	35:BA:450:G:H5'	2.43	0.48
35:BA:503:A:C6	35:BA:505:A:C6	3.02	0.48
35:BA:578:A:OP1	35:BA:1255:U:O2'	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:67:U:H2'	35:BA:68:G:C8	2.48	0.48
36:BB:90:A:C8	36:BB:91:C:H1'	2.48	0.48
37:BC:74:VAL:H	37:BC:91:ALA:HB1	1.77	0.48
35:BA:2631:G:H22	39:BE:61:ARG:NH1	2.10	0.48
40:BF:28:ILE:HA	40:BF:112:MET:HG2	1.94	0.48
40:BF:68:LYS:O	40:BF:69:HIS:CB	2.61	0.48
41:BG:36:LYS:HE3	41:BG:160:VAL:HG21	1.94	0.48
44:BN:107:LEU:HB2	44:BN:108:PRO:CD	2.41	0.48
45:BO:13:ASN:HD22	45:BO:97:ARG:HG2	1.78	0.48
45:BO:9:GLU:O	45:BO:83:ALA:HB1	2.13	0.48
46:BP:71:VAL:HG22	46:BP:72:PRO:HG3	1.94	0.48
48:BR:99:LYS:C	48:BR:100:LEU:HD22	2.34	0.48
49:BS:92:TYR:HD2	49:BS:97:ARG:NH1	2.11	0.48
51:BU:90:VAL:O	51:BU:92:ARG:N	2.44	0.48
51:BU:96:ALA:C	51:BU:98:LEU:H	2.17	0.48
52:BV:75:PHE:HB2	52:BV:87:HIS:CB	2.43	0.48
53:BW:64:MET:CE	53:BW:109:GLU:HG3	2.43	0.48
53:BW:64:MET:O	53:BW:65:LEU:CB	2.61	0.48
53:BW:6:ILE:HG22	53:BW:6:ILE:O	2.13	0.48
55:BY:60:PHE:HD2	55:BY:60:PHE:O	1.95	0.48
55:BY:14:LEU:O	55:BY:72:VAL:HG12	2.13	0.48
1:CA:112:G:H5'	1:CA:389:A:H4'	1.95	0.48
1:CA:113:G:H2'	1:CA:114:U:C6	2.49	0.48
1:CA:15:G:O3'	5:CE:24:ARG:NH2	2.46	0.48
1:CA:645:C:H2'	1:CA:646:U:H6	1.77	0.48
2:CB:17:PHE:CD2	2:CB:17:PHE:C	2.84	0.48
4:CD:131:ARG:HD3	4:CD:131:ARG:N	2.29	0.48
4:CD:159:ARG:HH11	4:CD:159:ARG:HG3	1.77	0.48
6:CF:8:ILE:HD13	6:CF:26:ILE:HD13	1.94	0.48
7:CG:100:ALA:O	7:CG:101:LEU:C	2.50	0.48
7:CG:122:HIS:O	7:CG:125:MET:HB3	2.13	0.48
8:CH:120:THR:O	8:CH:122:ARG:N	2.46	0.48
8:CH:28:ALA:HB2	8:CH:57:PRO:O	2.13	0.48
9:CI:56:LEU:HD23	9:CI:56:LEU:O	2.13	0.48
14:CN:4:LYS:HA	14:CN:7:ILE:HD11	1.94	0.48
18:CR:53:ARG:HH12	18:CR:59:SER:CA	2.22	0.48
26:D0:32:ARG:N	26:D0:35:ASN:HD22	1.93	0.48
28:D2:29:LYS:HA	28:D2:32:LEU:CB	2.39	0.48
35:DA:1127:A:H2'	35:DA:1128:A:H5''	1.94	0.48
35:DA:1184:G:C2'	35:DA:1185:C:H5'	2.43	0.48
35:DA:1798:U:N3	35:DA:1819:A:C2	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1992:G:N1	35:DA:1997:G:N1	2.60	0.48
35:DA:2065:C:H2'	35:DA:2066:C:H6	1.78	0.48
35:DA:2194:G:H5'	35:DA:2195:C:OP2	2.13	0.48
35:DA:2220:G:H2'	35:DA:2221:G:C8	2.36	0.48
35:DA:2290:G:H2'	35:DA:2291:U:O4'	2.13	0.48
35:DA:2336:A:H3'	35:DA:2337:G:H8	1.77	0.48
35:DA:238:C:H2'	35:DA:239:U:H6	1.78	0.48
35:DA:2431:U:H2'	35:DA:2433:A:OP2	2.13	0.48
35:DA:2695:C:H2'	35:DA:2696:U:C6	2.48	0.48
35:DA:2762:G:H2'	35:DA:2763:G:C5'	2.40	0.48
35:DA:828:U:H3'	35:DA:828:U:O2	2.12	0.48
37:DC:41:VAL:HB	37:DC:178:ALA:CB	2.43	0.48
38:DD:210:GLY:O	38:DD:211:ARG:CB	2.58	0.48
38:DD:211:ARG:O	38:DD:214:TRP:N	2.47	0.48
40:DF:110:LEU:HA	40:DF:183:VAL:CG1	2.39	0.48
41:DG:142:PRO:HG2	41:DG:143:GLU:HG2	1.95	0.48
41:DG:14:GLU:O	41:DG:17:PRO:HD2	2.13	0.48
42:DH:160:LYS:O	42:DH:163:TYR:HE1	1.96	0.48
35:DA:2726:U:H6	45:DO:67:LYS:NZ	2.11	0.48
46:DP:47:ASP:HB2	46:DP:51:PHE:CB	2.32	0.48
35:DA:244:A:O2'	46:DP:73:GLY:HA3	2.13	0.48
47:DQ:16:ARG:HB3	47:DQ:16:ARG:HH11	1.75	0.48
49:DS:89:ARG:CA	49:DS:89:ARG:NE	2.74	0.48
51:DU:108:GLU:OE1	51:DU:112:ARG:NH2	2.43	0.48
52:DV:45:THR:O	52:DV:45:THR:HG22	2.13	0.48
1:AA:1163:C:N3	1:AA:1174:G:N2	2.60	0.48
1:AA:1348:U:H2'	1:AA:1349:A:C8	2.49	0.48
1:AA:1519:A:N7	1:AA:1520:G:H1'	2.27	0.48
1:AA:262:A:C6	1:AA:263:A:C6	3.01	0.48
1:AA:241:C:H1'	1:AA:286:G:N2	2.28	0.48
1:AA:376:G:O2'	1:AA:377:G:H5'	2.14	0.48
2:AB:157:ARG:CG	2:AB:158:LEU:H	2.25	0.48
2:AB:24:TRP:O	2:AB:25:ASN:HB2	2.12	0.48
2:AB:68:ILE:HG22	2:AB:70:PHE:CD1	2.47	0.48
3:AC:153:VAL:HG12	3:AC:154:SER:H	1.78	0.48
3:AC:159:GLY:O	3:AC:160:ALA:C	2.51	0.48
4:AD:120:LEU:HB2	4:AD:126:ILE:HD11	1.94	0.48
4:AD:149:ALA:O	4:AD:150:GLU:C	2.51	0.48
5:AE:135:THR:O	5:AE:139:LEU:HG	2.14	0.48
8:AH:15:ASN:O	8:AH:16:ALA:C	2.52	0.48
11:AK:66:LEU:O	11:AK:67:ASP:C	2.52	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AL:89:ARG:NE	12:AL:91:LYS:HE2	2.28	0.48
13:AM:67:GLU:O	13:AM:68:GLY:C	2.52	0.48
14:AN:8:GLU:C	14:AN:12:ARG:HD3	2.34	0.48
17:AQ:34:LYS:O	17:AQ:36:ILE:HG12	2.13	0.48
17:AQ:45:HIS:O	17:AQ:73:VAL:HG23	2.12	0.48
25:AY:143:LEU:CD2	25:AY:144:ALA:N	2.77	0.48
27:B1:86:SER:O	27:B1:86:SER:OG	2.29	0.48
28:B2:60:LEU:O	28:B2:62:THR:N	2.46	0.48
32:B6:32:ASN:O	32:B6:33:LYS:HB2	2.13	0.48
35:BA:1186:G:H2'	35:BA:1187:G:H5'	1.94	0.48
35:BA:1424:G:O2'	35:BA:1425:G:H5'	2.13	0.48
35:BA:1591:G:C2'	35:BA:1592:C:H5'	2.43	0.48
35:BA:1678:G:OP2	35:BA:1678:G:C8	2.66	0.48
35:BA:1759:A:H2'	35:BA:1760:A:H8	1.76	0.48
35:BA:1835:G:N3	35:BA:1835:G:H2'	2.28	0.48
35:BA:2388:A:H5'	35:BA:2389:G:OP2	2.13	0.48
35:BA:2666:C:O2'	35:BA:2667:C:H5'	2.12	0.48
35:BA:407:G:H2'	35:BA:408:G:C8	2.46	0.48
35:BA:660:G:C6	35:BA:661:C:C4	3.01	0.48
35:BA:666:G:O2'	35:BA:667:U:H5'	2.12	0.48
35:BA:998:C:P	51:BU:93:LYS:HE2	2.53	0.48
36:BB:18:G:H2'	36:BB:19:G:H8	1.78	0.48
36:BB:28:C:O2'	36:BB:29:A:H5'	2.13	0.48
36:BB:52:A:O2'	36:BB:53:A:H8	1.96	0.48
38:BD:132:PRO:HG3	38:BD:190:TYR:CE1	2.48	0.48
39:BE:14:ILE:HB	50:BT:14:TYR:HE2	1.79	0.48
39:BE:4:ILE:HD11	39:BE:29:GLY:H	1.78	0.48
39:BE:38:THR:CG2	39:BE:39:PRO:HD2	2.44	0.48
39:BE:60:ASN:O	39:BE:62:PRO:N	2.47	0.48
40:BF:70:THR:O	40:BF:72:ARG:N	2.46	0.48
40:BF:74:ARG:O	40:BF:75:HIS:ND1	2.45	0.48
41:BG:120:LEU:HB3	41:BG:131:TYR:OH	2.13	0.48
41:BG:45:GLU:HG2	41:BG:45:GLU:O	2.13	0.48
43:BI:10:GLU:OE1	43:BI:11:ASN:N	2.46	0.48
48:BR:26:LYS:HG2	48:BR:70:LEU:HD22	1.95	0.48
49:BS:24:LEU:H	49:BS:24:LEU:HD22	1.78	0.48
50:BT:32:TYR:HB2	50:BT:33:LYS:H	1.46	0.48
51:BU:20:LEU:H	51:BU:20:LEU:CD2	2.21	0.48
52:BV:36:PRO:HD2	52:BV:60:GLU:O	2.14	0.48
52:BV:75:PHE:HB2	52:BV:87:HIS:HB3	1.95	0.48
53:BW:45:TYR:CD2	53:BW:45:TYR:C	2.86	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:86:LEU:HD12	53:BW:87:PRO:HD2	1.95	0.48
54:BX:72:LYS:O	54:BX:73:ARG:HB3	2.13	0.48
56:BZ:10:ARG:CZ	56:BZ:36:LYS:HB3	2.43	0.48
56:BZ:52:SER:OG	56:BZ:53:ILE:N	2.46	0.48
1:CA:1001:A:H2	1:CA:1001(A):G:O6	1.96	0.48
1:CA:1178:G:N2	1:CA:1180:A:H3'	2.28	0.48
1:CA:542:G:C6	1:CA:543:C:N4	2.82	0.48
1:CA:955:U:O2'	1:CA:956:U:H5'	2.13	0.48
4:CD:170:VAL:CG1	4:CD:174:LEU:HB2	2.43	0.48
5:CE:26:PHE:O	5:CE:27:ARG:HB2	2.12	0.48
7:CG:40:ALA:O	7:CG:44:TYR:HD1	1.96	0.48
13:CM:67:GLU:O	13:CM:68:GLY:C	2.52	0.48
13:CM:91:ARG:O	13:CM:96:LEU:N	2.46	0.48
15:CO:33:THR:HG23	15:CO:63:ARG:HH11	1.78	0.48
16:CP:6:LEU:HD23	16:CP:17:TYR:CD2	2.48	0.48
18:CR:19:LYS:O	18:CR:20:ALA:CB	2.61	0.48
18:CR:73:ALA:HB3	18:CR:79:LEU:HD12	1.96	0.48
25:CY:122:ALA:O	25:CY:125:GLY:N	2.45	0.48
27:D1:23:LYS:CB	27:D1:23:LYS:HZ2	2.25	0.48
29:D3:41:PRO:HB3	35:DA:852:G:O2'	2.13	0.48
32:D6:13:CYS:O	32:D6:21:TYR:HA	2.13	0.48
35:DA:132:G:H2'	35:DA:133:C:C6	2.48	0.48
35:DA:134:C:O2'	35:DA:135:G:H5'	2.13	0.48
35:DA:1514:U:C2	35:DA:1515:G:C8	3.02	0.48
35:DA:1563:G:O2'	35:DA:1564:C:H5'	2.13	0.48
35:DA:1268:A:C6	35:DA:2013:A:C8	3.02	0.48
35:DA:2364:C:C2'	35:DA:2365:G:H5'	2.43	0.48
35:DA:2814:C:H2'	35:DA:2815:C:C6	2.45	0.48
35:DA:2841:C:C2	35:DA:2877:G:C2	3.01	0.48
35:DA:2882:A:OP1	48:DR:96:ARG:HD3	2.12	0.48
35:DA:660:G:C6	35:DA:661:C:C4	3.01	0.48
35:DA:785:G:H2'	35:DA:786:C:C6	2.48	0.48
35:DA:790:C:O2'	35:DA:791:C:C5'	2.61	0.48
35:DA:830:G:C4	35:DA:2448:A:C5	3.01	0.48
36:DB:25:A:H2'	36:DB:26:A:O4'	2.13	0.48
36:DB:65:C:O2'	36:DB:66:A:H5'	2.12	0.48
39:DE:161:GLY:O	39:DE:163:GLU:HG2	2.13	0.48
40:DF:157:VAL:HB	40:DF:194:MET:CB	2.43	0.48
40:DF:28:ILE:CD1	40:DF:28:ILE:H	2.26	0.48
35:DA:797:C:OP2	40:DF:62:ARG:HG3	2.13	0.48
42:DH:65:HIS:CG	42:DH:66:GLY:N	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DI:112:LYS:H	43:DI:114:LEU:HG	1.79	0.48
43:DI:145:VAL:CG1	43:DI:146:ALA:N	2.76	0.48
47:DQ:16:ARG:HG2	47:DQ:17:LEU:N	2.21	0.48
49:DS:15:ARG:C	49:DS:17:ARG:N	2.65	0.48
52:DV:61:VAL:CG2	52:DV:99:ILE:HB	2.39	0.48
53:DW:76:VAL:HB	53:DW:103:ILE:HG22	1.95	0.48
53:DW:64:MET:O	53:DW:65:LEU:CB	2.61	0.48
55:DY:74:PRO:HG2	55:DY:80:GLY:C	2.34	0.48
56:DZ:49:ARG:HB2	56:DZ:50:GLN:OE1	2.13	0.48
1:AA:1236:A:H4'	1:AA:1304:G:H4'	1.95	0.48
1:AA:1422:G:C2	1:AA:1479:C:N3	2.81	0.48
1:AA:1480:G:N1	1:AA:1481:U:C2	2.81	0.48
1:AA:189(D):C:H1'	1:AA:189(H):G:N2	2.28	0.48
1:AA:193:C:O2'	1:AA:194:C:H5'	2.14	0.48
1:AA:391:G:C6	1:AA:392:G:N7	2.81	0.48
1:AA:66:G:N3	1:AA:66:G:H2'	2.28	0.48
1:AA:81:U:H2'	1:AA:82:U:C6	2.48	0.48
1:AA:955:U:O2'	1:AA:956:U:H5'	2.13	0.48
3:AC:7:PRO:O	3:AC:11:ARG:HG2	2.13	0.48
4:AD:154:ASN:HA	4:AD:159:ARG:HE	1.78	0.48
4:AD:158:ILE:O	4:AD:162:LEU:HG	2.13	0.48
4:AD:56:VAL:C	4:AD:58:LEU:N	2.66	0.48
7:AG:115:ARG:O	7:AG:118:VAL:HG23	2.13	0.48
7:AG:119:ARG:O	7:AG:120:ILE:C	2.52	0.48
25:AY:60:ALA:HB1	25:AY:65:THR:O	2.13	0.48
26:B0:27:GLU:H	26:B0:69:PHE:HE1	1.58	0.48
27:B1:89:GLU:C	27:B1:90:ILE:HD13	2.33	0.48
33:B7:8:ASN:ND2	33:B7:8:ASN:C	2.66	0.48
34:B8:4:MET:CE	34:B8:61:LEU:HD12	2.44	0.48
35:BA:1247:A:OP1	40:BF:95:ARG:NH2	2.47	0.48
35:BA:1360:A:C5	35:BA:1372:U:C4	3.01	0.48
35:BA:1688:U:O2	35:BA:1688:U:H2'	2.12	0.48
35:BA:1795:C:N4	35:BA:1824:G:H1	2.11	0.48
35:BA:1927:A:C6	35:BA:1928:A:C6	3.02	0.48
35:BA:1991:U:C2'	35:BA:1992:G:C5'	2.91	0.48
35:BA:2052:G:H4'	39:BE:143:ASN:O	2.13	0.48
35:BA:2460:U:C2'	35:BA:2461:C:H5'	2.44	0.48
35:BA:2643:G:C2	35:BA:2772:C:N3	2.81	0.48
35:BA:2696:U:H2'	35:BA:2697:G:H8	1.78	0.48
35:BA:2839:G:H1'	48:BR:93:GLY:H	1.77	0.48
35:BA:580:C:H2'	35:BA:581:C:H6	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:681:G:H2'	35:BA:682:G:O4'	2.12	0.48
36:BB:5:C:H2'	36:BB:6:C:H6	1.78	0.48
38:BD:139:GLY:H	38:BD:165:ILE:HB	1.79	0.48
38:BD:44:ASN:ND2	38:BD:47:GLY:O	2.46	0.48
39:BE:116:VAL:CG2	39:BE:122:PHE:HB2	2.38	0.48
41:BG:76:SER:HB3	41:BG:84:LYS:N	2.21	0.48
42:BH:160:LYS:O	42:BH:163:TYR:HE1	1.95	0.48
43:BI:83:ALA:HB2	43:BI:88:ILE:HG23	1.94	0.48
44:BN:71:ILE:HD13	44:BN:86:PRO:HA	1.94	0.48
45:BO:68:GLU:HB3	45:BO:78:ARG:CD	2.41	0.48
47:BQ:39:PRO:O	47:BQ:40:ALA:HB2	2.13	0.48
48:BR:28:LEU:O	48:BR:30:THR:N	2.45	0.48
35:BA:2378:A:O2'	49:BS:22:GLY:HA3	2.13	0.48
49:BS:90:GLY:C	49:BS:92:TYR:H	2.17	0.48
51:BU:108:GLU:HB3	51:BU:112:ARG:NH1	2.27	0.48
51:BU:86:ALA:HB2	51:BU:116:ALA:HB1	1.95	0.48
52:BV:18:LEU:HA	52:BV:97:LYS:HZ1	1.78	0.48
53:BW:70:TYR:CE2	53:BW:108:GLY:HA3	2.42	0.48
55:BY:44:ILE:HG23	55:BY:45:VAL:N	2.28	0.48
55:BY:88:LYS:NZ	55:BY:93:GLY:N	2.50	0.48
56:BZ:76:LEU:HB2	56:BZ:81:ARG:O	2.14	0.48
56:BZ:27:VAL:N	56:BZ:85:HIS:HE1	2.11	0.48
1:CA:1042:G:O2'	1:CA:1043:C:H5'	2.14	0.48
1:CA:1104:G:H2'	1:CA:1105:A:H8	1.77	0.48
1:CA:1473:A:O2'	1:CA:1474:G:H5'	2.14	0.48
1:CA:262:A:C6	1:CA:263:A:C6	3.02	0.48
1:CA:314:C:O2'	1:CA:315:A:H5'	2.12	0.48
1:CA:541:G:H2'	1:CA:542:G:H8	1.79	0.48
1:CA:601:C:H42	1:CA:637:G:H1	1.62	0.48
3:CC:7:PRO:O	3:CC:11:ARG:HG2	2.13	0.48
3:CC:92:ALA:C	3:CC:94:LEU:H	2.17	0.48
4:CD:176:LEU:HD12	4:CD:177:ASP:H	1.78	0.48
5:CE:149:GLU:C	5:CE:151:LEU:N	2.66	0.48
5:CE:8:GLU:N	5:CE:34:VAL:HG13	2.28	0.48
6:CF:3:ARG:NH1	6:CF:38:GLU:OE2	2.46	0.48
6:CF:69:GLU:O	6:CF:70:ASP:C	2.52	0.48
7:CG:69:VAL:O	7:CG:69:VAL:HG12	2.12	0.48
7:CG:79:ARG:HH11	7:CG:79:ARG:HG3	1.77	0.48
8:CH:6:ILE:CG2	8:CH:10:LEU:HD11	2.42	0.48
12:CL:85:ILE:CD1	12:CL:98:TYR:HB3	2.42	0.48
16:CP:48:TRP:O	16:CP:49:LEU:C	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:63:GLN:OE1	18:CR:63:GLN:HA	2.13	0.48
18:CR:74:ARG:HE	18:CR:81:PHE:CA	2.22	0.48
20:CT:38:LYS:O	20:CT:39:LYS:C	2.51	0.48
21:CU:21:TYR:N	21:CU:21:TYR:CD1	2.80	0.48
25:CY:64:ARG:HA	25:CY:103:ILE:CG1	2.44	0.48
25:CY:92:PRO:HB3	25:CY:101:ILE:HG12	1.95	0.48
27:D1:49:VAL:O	27:D1:62:VAL:O	2.31	0.48
27:D1:90:ILE:O	27:D1:93:GLU:HG2	2.13	0.48
30:D4:45:GLY:C	30:D4:47:GLN:N	2.66	0.48
35:DA:1230:C:O2'	35:DA:1231:G:H5'	2.13	0.48
35:DA:1369:G:N2	35:DA:1370:C:C2	2.82	0.48
35:DA:1400:G:H2'	35:DA:1401:G:C8	2.48	0.48
35:DA:1785:A:N7	35:DA:1787:A:C5	2.81	0.48
27:D1:41:ARG:NH2	35:DA:205:G:N1	2.58	0.48
35:DA:2165:G:H3'	35:DA:2166:G:H8	1.78	0.48
35:DA:2170:A:H5'	35:DA:2171:A:OP2	2.13	0.48
35:DA:2250:G:C6	47:DQ:82:ARG:HD2	2.48	0.48
35:DA:2711:A:OP1	35:DA:2712(A):A:OP1	2.32	0.48
35:DA:37:C:H2'	35:DA:38:A:H8	1.78	0.48
35:DA:543:C:C4	35:DA:551:G:N1	2.80	0.48
35:DA:588:U:H1'	40:DF:90:PHE:HB3	1.95	0.48
35:DA:660:G:C5'	40:DF:99:TYR:CD2	2.95	0.48
35:DA:662:G:O2'	35:DA:663:G:H5'	2.13	0.48
36:DB:66:A:N6	36:DB:108:U:H2'	2.25	0.48
39:DE:117:MET:CE	39:DE:124:GLY:HA3	2.44	0.48
39:DE:23:VAL:HA	39:DE:184:VAL:O	2.14	0.48
40:DF:171:PRO:HG2	40:DF:172:TRP:H	1.79	0.48
45:DO:65:THR:O	45:DO:79:PHE:HB2	2.13	0.48
48:DR:94:TYR:CD1	48:DR:94:TYR:N	2.82	0.48
51:DU:86:ALA:HB2	51:DU:116:ALA:HB1	1.94	0.48
54:DX:57:LEU:HD12	54:DX:76:ARG:CD	2.42	0.48
56:DZ:121:HIS:C	56:DZ:123:ASP:N	2.66	0.48
56:DZ:140:ASP:C	56:DZ:142:SER:H	2.16	0.48
1:AA:1152:A:H2'	1:AA:1153:C:C6	2.47	0.48
1:AA:1166:G:H2'	1:AA:1169:A:OP2	2.14	0.48
1:AA:1234:C:C4'	1:AA:1364:U:H1'	2.43	0.48
1:AA:1238:A:H62	1:AA:1301:U:H3	1.62	0.48
1:AA:1349:A:O2'	1:AA:1350:A:H5'	2.13	0.48
1:AA:66:G:C4'	1:AA:173:U:C5	2.94	0.48
1:AA:332:G:H2'	1:AA:333:G:C8	2.43	0.48
1:AA:438:G:H4'	4:AD:123:HIS:ND1	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:473:G:H2'	1:AA:474:G:C8	2.47	0.48
1:AA:615:C:H2'	1:AA:616:G:O4'	2.12	0.48
1:AA:640:A:H4'	8:AH:116:LYS:HZ1	1.79	0.48
1:AA:666:G:H1'	1:AA:741:G:H22	1.79	0.48
1:AA:893:C:H2'	1:AA:894:G:H8	1.75	0.48
2:AB:95:GLN:HB3	2:AB:148:TYR:HD1	1.78	0.48
2:AB:42:ILE:CD1	2:AB:203:GLY:H	2.24	0.48
3:AC:178:LEU:N	3:AC:178:LEU:HD22	2.29	0.48
4:AD:13:ARG:CG	4:AD:14:ARG:N	2.75	0.48
4:AD:15:GLU:CD	4:AD:15:GLU:H	2.17	0.48
6:AF:16:GLN:O	6:AF:19:LEU:N	2.47	0.48
7:AG:146:GLU:CA	7:AG:149:ARG:HB2	2.44	0.48
7:AG:47:CYS:HB3	7:AG:58:PRO:CG	2.44	0.48
7:AG:149:ARG:HD3	11:AK:59:TYR:CZ	2.48	0.48
12:AL:41:ARG:CG	12:AL:42:THR:H	2.26	0.48
12:AL:58:VAL:O	12:AL:60:LEU:HD22	2.14	0.48
17:AQ:11:VAL:CA	17:AQ:53:LEU:HD11	2.43	0.48
27:B1:87:PRO:HD2	27:B1:89:GLU:OE1	2.13	0.48
31:B5:32:PRO:O	31:B5:33:CYS:CB	2.61	0.48
34:B8:56:GLU:HA	34:B8:59:LYS:HZ1	1.78	0.48
35:BA:1024:G:P	35:BA:1025:G:H3'	2.53	0.48
35:BA:1500:G:C6	35:BA:1501:C:N3	2.82	0.48
35:BA:1902:C:H2'	35:BA:1903:G:O4'	2.13	0.48
35:BA:2175:C:C3'	35:BA:2176:A:H5''	2.43	0.48
35:BA:2306:C:H5'	35:BA:2307:G:O5'	2.13	0.48
34:B8:30:ARG:HG2	35:BA:2393:A:H5''	1.95	0.48
35:BA:2460:U:H2'	35:BA:2461:C:O4'	2.13	0.48
35:BA:2692:C:N4	35:BA:2718:G:N1	2.62	0.48
35:BA:2831:G:O4'	35:BA:2883:A:C2	2.66	0.48
35:BA:314:A:O2'	35:BA:315:G:H5'	2.13	0.48
35:BA:538:G:N3	35:BA:538:G:H2'	2.29	0.48
35:BA:998:C:H42	35:BA:1157:G:H1	1.60	0.48
37:BC:83:ILE:O	37:BC:83:ILE:HG22	2.14	0.48
38:BD:113:VAL:C	38:BD:115:GLN:H	2.15	0.48
38:BD:118:VAL:HG22	38:BD:119:ALA:H	1.77	0.48
38:BD:27:THR:O	38:BD:29:PRO:HD2	2.12	0.48
38:BD:25:THR:HG21	38:BD:82:ILE:H	1.72	0.48
39:BE:199:ARG:HG3	39:BE:199:ARG:NH1	2.28	0.48
40:BF:125:LEU:HD12	40:BF:196:LEU:HD22	1.95	0.48
40:BF:114:VAL:CG1	40:BF:202:PHE:HE2	2.25	0.48
40:BF:28:ILE:CD1	40:BF:28:ILE:H	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:63:LYS:HZ3	40:BF:67:GLN:HB3	1.75	0.48
41:BG:91:ARG:HD2	41:BG:92:VAL:N	2.27	0.48
42:BH:123:PHE:HA	42:BH:133:VAL:CG2	2.36	0.48
44:BN:128:HIS:C	44:BN:130:HIS:HD1	2.17	0.48
44:BN:38:HIS:CG	44:BN:39:ARG:H	2.31	0.48
44:BN:43:THR:O	44:BN:46:VAL:HG12	2.14	0.48
46:BP:41:ARG:CD	46:BP:41:ARG:N	2.76	0.48
47:BQ:118:LEU:O	47:BQ:121:ALA:N	2.46	0.48
47:BQ:141:GLN:NE2	56:BZ:89:PHE:HB3	2.28	0.48
47:BQ:42:ILE:N	47:BQ:42:ILE:HD12	2.28	0.48
48:BR:18:LEU:HD13	48:BR:18:LEU:C	2.33	0.48
48:BR:39:PRO:O	48:BR:40:LYS:C	2.52	0.48
48:BR:53:HIS:O	48:BR:56:LYS:CB	2.62	0.48
50:BT:80:SER:CB	50:BT:81:PRO:HD2	2.35	0.48
50:BT:96:ARG:CZ	50:BT:96:ARG:HB3	2.42	0.48
53:BW:12:ILE:HD12	53:BW:42:ARG:HH11	1.78	0.48
54:BX:84:ALA:O	54:BX:86:GLY:N	2.47	0.48
1:CA:1163:C:N3	1:CA:1174:G:N2	2.61	0.48
1:CA:1289:A:H3'	1:CA:1290:G:H8	1.79	0.48
1:CA:1348:U:H2'	1:CA:1349:A:C8	2.49	0.48
1:CA:389:A:H2'	1:CA:390:C:C5'	2.43	0.48
2:CB:95:GLN:HB3	2:CB:148:TYR:HD1	1.77	0.48
2:CB:80:ILE:HG12	2:CB:211:ILE:CG2	2.43	0.48
7:CG:102:ARG:O	7:CG:103:TRP:C	2.50	0.48
8:CH:65:TYR:HA	8:CH:79:VAL:HG23	1.95	0.48
10:CJ:56:HIS:O	10:CJ:58:ASP:O	2.32	0.48
11:CK:122:LYS:O	11:CK:126:ARG:HB3	2.14	0.48
15:CO:56:LEU:HA	15:CO:59:MET:CE	2.43	0.48
19:CS:36:ARG:CZ	19:CS:75:ALA:HB3	2.44	0.48
25:CY:135:GLU:O	25:CY:138:ASP:HB2	2.13	0.48
25:CY:139:LYS:O	25:CY:140:LEU:C	2.51	0.48
25:CY:163:LYS:O	25:CY:164:ILE:C	2.52	0.48
25:CY:30:THR:HB	25:CY:182:GLU:OE2	2.14	0.48
25:CY:80:GLU:HG3	25:CY:94:ASN:OD1	2.14	0.48
27:D1:15:ALA:O	27:D1:46:LEU:CD2	2.61	0.48
27:D1:93:GLU:N	27:D1:93:GLU:OE2	2.47	0.48
28:D2:14:ARG:HE	28:D2:14:ARG:CA	2.27	0.48
32:D6:47:THR:HG22	32:D6:49:HIS:H	1.77	0.48
32:D6:51:GLU:CG	32:D6:52:VAL:H	2.00	0.48
33:D7:41:ARG:O	33:D7:42:LEU:C	2.52	0.48
35:DA:1196:C:H2'	35:DA:1197:G:H8	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1260:G:H2'	35:DA:1261:C:C6	2.49	0.48
35:DA:1305:C:O2'	35:DA:1306:C:H5'	2.14	0.48
35:DA:1367:A:H3'	35:DA:1368:G:O4'	2.13	0.48
35:DA:1543:C:O2	35:DA:1543:C:C2'	2.60	0.48
31:D5:19:ARG:HG3	35:DA:2046:G:H5'	1.95	0.48
35:DA:2505:G:H2'	35:DA:2576:G:C6	2.49	0.48
35:DA:2642:G:N2	35:DA:2773:C:C2	2.81	0.48
35:DA:2677:G:C4	35:DA:2678:C:H5	2.31	0.48
35:DA:2689:U:H4'	35:DA:2690:C:H6	1.77	0.48
35:DA:2736:G:O2'	35:DA:2737:G:H5'	2.13	0.48
35:DA:444:C:O2'	35:DA:445:C:H5'	2.13	0.48
35:DA:469:G:O2'	35:DA:470:A:H5''	2.14	0.48
35:DA:481:G:C2'	35:DA:482:A:OP2	2.61	0.48
35:DA:733:G:O5'	35:DA:733:G:H8	1.96	0.48
35:DA:745:G:C2'	35:DA:746:A:H5'	2.44	0.48
35:DA:802:A:N3	35:DA:802:A:H2'	2.28	0.48
35:DA:869:G:O2'	47:DQ:8:LYS:HD3	2.13	0.48
35:DA:869:G:H2'	35:DA:870:A:C8	2.48	0.48
35:DA:917:A:H2'	35:DA:918:A:O4'	2.13	0.48
35:DA:951:C:O2'	35:DA:952:G:H5'	2.12	0.48
35:DA:998:C:N4	35:DA:1158:C:H42	2.12	0.48
36:DB:115:G:H2'	36:DB:116:G:H8	1.78	0.48
45:DO:93:PRO:C	45:DO:95:GLY:N	2.67	0.48
47:DQ:118:LEU:O	47:DQ:121:ALA:N	2.47	0.48
47:DQ:41:TRP:HB3	47:DQ:94:VAL:HG21	1.95	0.48
48:DR:117:VAL:O	48:DR:118:GLU:HB2	2.13	0.48
35:DA:2378:A:O2'	49:DS:22:GLY:HA3	2.12	0.48
49:DS:79:ALA:C	49:DS:80:LEU:HD12	2.33	0.48
49:DS:89:ARG:HB3	49:DS:97:ARG:HH22	1.79	0.48
45:DO:104:ARG:NH1	50:DT:35:LYS:HB3	2.28	0.48
54:DX:47:PHE:N	54:DX:47:PHE:CD1	2.81	0.48
56:DZ:30:ASN:C	56:DZ:32:HIS:N	2.65	0.48
56:DZ:3:TYR:N	56:DZ:56:VAL:O	2.46	0.48
1:AA:1507:A:H2	1:AA:1530:G:C1'	2.24	0.48
1:AA:439:A:C4	1:AA:496:A:C2	3.01	0.48
1:AA:516:U:O2'	1:AA:517:G:H5'	2.14	0.48
1:AA:880:C:H2'	1:AA:881:G:H8	1.77	0.48
3:AC:116:VAL:O	3:AC:119:ARG:HB3	2.14	0.48
3:AC:40:ARG:O	3:AC:41:GLY:C	2.51	0.48
3:AC:84:ILE:HA	3:AC:87:LEU:HB2	1.96	0.48
3:AC:91:LEU:CB	3:AC:99:VAL:HG21	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:13:ARG:O	4:AD:15:GLU:N	2.47	0.48
4:AD:68:TYR:CZ	4:AD:97:LEU:HD22	2.48	0.48
6:AF:67:MET:HE2	6:AF:72:VAL:H	1.78	0.48
1:AA:1343:G:H4'	9:AI:122:ALA:O	2.13	0.48
13:AM:91:ARG:O	13:AM:96:LEU:N	2.47	0.48
19:AS:16:LEU:O	19:AS:20:LEU:N	2.44	0.48
27:B1:47:GLN:O	27:B1:64:ALA:HB2	2.14	0.48
29:B3:59:VAL:CG1	29:B3:60:GLU:N	2.71	0.48
34:B8:29:LYS:NZ	34:B8:44:LYS:CB	2.77	0.48
35:BA:1006:C:H2'	35:BA:1007:C:H6	1.77	0.48
35:BA:111:A:O2'	35:BA:112:U:H5'	2.13	0.48
35:BA:2323:G:H2'	35:BA:2324:C:O4'	2.12	0.48
35:BA:2481:G:HO2'	35:BA:2482:G:P	2.35	0.48
35:BA:2639:A:C2'	35:BA:2640:G:C5'	2.85	0.48
35:BA:494:G:O2'	35:BA:495:G:H5'	2.13	0.48
35:BA:514:A:H2'	35:BA:515:A:H8	1.77	0.48
31:B5:3:LYS:HB2	35:BA:747:U:C5	2.48	0.48
35:BA:778:G:H4'	38:BD:48:ARG:HD2	1.96	0.48
35:BA:674:G:H2'	35:BA:804:A:H61	1.77	0.48
35:BA:1566:A:C4	38:BD:214:TRP:CE3	3.02	0.48
39:BE:31:CYS:HB3	39:BE:49:LEU:HB3	1.95	0.48
40:BF:39:TRP:CE3	40:BF:40:GLN:HG2	2.48	0.48
41:BG:105:LYS:HE2	41:BG:143:GLU:OE2	2.14	0.48
41:BG:19:LEU:HD21	41:BG:175:LEU:HD11	1.94	0.48
41:BG:34:LEU:HD12	41:BG:100:TRP:CZ3	2.48	0.48
41:BG:73:ALA:O	41:BG:85:GLY:HA2	2.13	0.48
45:BO:104:ARG:O	45:BO:107:ARG:N	2.42	0.48
45:BO:11:ALA:HB2	45:BO:64:ARG:NH2	2.28	0.48
46:BP:38:GLN:O	46:BP:39:LYS:HB2	2.13	0.48
47:BQ:87:LYS:O	47:BQ:87:LYS:CG	2.52	0.48
48:BR:63:ARG:C	48:BR:67:LEU:HD23	2.34	0.48
50:BT:109:GLU:HA	50:BT:112:ARG:HD2	1.95	0.48
54:BX:83:VAL:C	54:BX:85:PRO:CD	2.82	0.48
1:CA:1054:C:P	1:CA:1197:G:OP2	2.71	0.48
1:CA:1260:C:H4'	1:CA:1284:C:H5'	1.96	0.48
1:CA:1279:A:H2'	1:CA:1279:A:N3	2.28	0.48
1:CA:1306:A:H2'	1:CA:1307:U:O4'	2.12	0.48
1:CA:1323:G:H2'	1:CA:1324:A:C8	2.48	0.48
1:CA:599:C:H2'	1:CA:600:C:C6	2.47	0.48
1:CA:615:C:H2'	1:CA:616:G:O4'	2.14	0.48
2:CB:145:LEU:HD22	2:CB:149:LEU:HD12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1057:G:H5''	3:CC:154:SER:HG	1.77	0.48
3:CC:173:VAL:N	3:CC:174:PRO:CD	2.77	0.48
3:CC:182:ILE:HG12	3:CC:203:PHE:HD1	1.79	0.48
4:CD:196:LEU:HB3	4:CD:197:PRO:HD2	1.94	0.48
4:CD:3:ARG:HD3	4:CD:3:ARG:O	2.13	0.48
5:CE:71:LEU:HD11	5:CE:114:GLY:C	2.33	0.48
6:CF:68:PRO:HG3	6:CF:71:ARG:HE	1.79	0.48
7:CG:92:SER:OG	7:CG:93:PRO:HD2	2.13	0.48
11:CK:34:ASP:HB2	11:CK:35:PRO:CD	2.43	0.48
12:CL:25:PRO:O	12:CL:27:LEU:HD22	2.14	0.48
13:CM:3:ARG:HG2	13:CM:9:ILE:HD11	1.94	0.48
15:CO:85:LEU:HD12	15:CO:87:ILE:HD11	1.94	0.48
19:CS:36:ARG:HH12	19:CS:75:ALA:CB	2.24	0.48
26:D0:56:ASP:CG	26:D0:58:THR:HG1	2.16	0.48
30:D4:14:ILE:CB	41:DG:5:VAL:HG22	2.44	0.48
33:D7:24:THR:CG2	33:D7:27:GLY:HA3	2.35	0.48
34:D8:39:LYS:O	34:D8:42:ARG:HB3	2.13	0.48
34:D8:39:LYS:HZ2	34:D8:40:GLU:HA	1.78	0.48
34:D8:61:LEU:O	34:D8:64:TYR:HD1	1.95	0.48
35:DA:105:C:C5'	35:DA:106:C:OP2	2.61	0.48
35:DA:1246:A:P	46:DP:16:ARG:HH12	2.36	0.48
35:DA:1360:A:C6	35:DA:1372:U:C5	3.02	0.48
35:DA:1472:A:O2'	35:DA:1473:G:H5'	2.14	0.48
35:DA:154:G:H1	35:DA:172:C:N4	2.09	0.48
35:DA:1598:C:H5'	54:DX:37:THR:HB	1.96	0.48
35:DA:1927:A:C6	35:DA:1928:A:C6	3.02	0.48
35:DA:2052:G:H4'	39:DE:143:ASN:O	2.14	0.48
35:DA:2239:G:C5	35:DA:2240:C:C5	3.02	0.48
35:DA:2552:U:C2	35:DA:2554:U:H5'	2.48	0.48
35:DA:271(V):G:C2	35:DA:271(W):G:H1'	2.48	0.48
35:DA:8:A:H2	35:DA:2896:C:O2	1.97	0.48
35:DA:460:A:C6	35:DA:470:A:C8	3.01	0.48
33:D7:16:HIS:ND1	35:DA:465:G:H4'	2.29	0.48
35:DA:576:U:H2'	35:DA:577:G:C8	2.49	0.48
35:DA:261:G:C1'	35:DA:609:A:H2	2.27	0.48
35:DA:911:A:H5''	35:DA:912:C:H5''	1.96	0.48
39:DE:16:ARG:O	39:DE:17:ASP:HB3	2.13	0.48
39:DE:8:LYS:HD2	39:DE:189:PRO:O	2.14	0.48
41:DG:49:ASP:O	41:DG:50:ALA:CB	2.62	0.48
41:DG:4:ASP:HB3	41:DG:8:LYS:HE3	1.94	0.48
42:DH:125:VAL:HG12	42:DH:127:GLU:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:89:ILE:HG12	42:DH:90:LYS:N	2.28	0.48
43:DI:111:PRO:HA	43:DI:114:LEU:HD11	1.94	0.48
43:DI:91:SER:H	43:DI:121:LYS:HE3	1.77	0.48
44:DN:67:LEU:HD22	44:DN:88:GLU:HG2	1.96	0.48
35:DA:833:U:P	46:DP:45:LEU:HD11	2.54	0.48
48:DR:52:ILE:HD12	48:DR:79:LEU:HD21	1.94	0.48
49:DS:28:VAL:N	49:DS:89:ARG:HG2	2.29	0.48
51:DU:108:GLU:HA	51:DU:111:GLU:HG2	1.95	0.48
53:DW:69:LEU:HA	53:DW:109:GLU:HA	1.95	0.48
53:DW:92:ARG:NH1	53:DW:92:ARG:HG2	2.29	0.48
56:DZ:134:PRO:O	56:DZ:135:GLU:C	2.52	0.48
1:AA:1237:C:H4'	1:AA:1334:G:N2	2.28	0.48
1:AA:1422:G:N2	1:AA:1479:C:N3	2.62	0.48
1:AA:1499:A:H2'	1:AA:1500:A:H5'	1.94	0.48
1:AA:329:A:C2	1:AA:332:G:C5	3.02	0.48
1:AA:336:C:H2'	1:AA:337:C:H6	1.79	0.48
1:AA:389:A:H2'	1:AA:390:C:C5'	2.43	0.48
1:AA:652:U:H2'	1:AA:752:G:N1	2.28	0.48
1:AA:705:U:C5	1:AA:706:A:C5	3.02	0.48
1:AA:575:G:N1	1:AA:821:G:C5	2.81	0.48
1:AA:919:A:C2'	1:AA:920:U:O5'	2.62	0.48
2:AB:193:ASP:O	2:AB:194:PRO:O	2.32	0.48
2:AB:39:ILE:HG22	2:AB:40:HIS:N	2.28	0.48
4:AD:154:ASN:CA	4:AD:159:ARG:HH21	2.27	0.48
4:AD:42:GLN:C	4:AD:42:GLN:OE1	2.52	0.48
4:AD:92:VAL:O	4:AD:96:LEU:HD22	2.14	0.48
5:AE:37:ARG:HA	5:AE:112:LEU:O	2.13	0.48
7:AG:104:LEU:CD2	7:AG:104:LEU:H	2.26	0.48
1:AA:1349:A:OP1	9:AI:120:ARG:HB3	2.13	0.48
12:AL:54:LYS:CD	12:AL:54:LYS:N	2.74	0.48
16:AP:8:ARG:NH1	16:AP:8:ARG:HG2	2.25	0.48
18:AR:71:LYS:O	18:AR:75:ILE:HG13	2.14	0.48
32:B6:52:VAL:HG12	32:B6:52:VAL:O	2.13	0.48
35:BA:1127:A:H2'	35:BA:1128:A:H5''	1.94	0.48
35:BA:1166:C:H2'	35:BA:1167:U:H6	1.78	0.48
35:BA:1233:C:C2'	35:BA:1234:U:H5'	2.43	0.48
35:BA:1281:G:H1	35:BA:1286:A:H62	1.62	0.48
35:BA:1309:G:C2'	35:BA:1310:G:H5'	2.44	0.48
35:BA:768:G:O2'	35:BA:1379:A:N6	2.45	0.48
35:BA:1516:C:H2'	35:BA:1517:G:C8	2.47	0.48
35:BA:1497:U:H3	35:BA:1578:U:P	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2006:C:H2'	35:BA:2007:C:C6	2.48	0.48
35:BA:2102:U:H2'	35:BA:2103:C:C5	2.48	0.48
35:BA:2189:U:H2'	35:BA:2190:G:O4'	2.14	0.48
35:BA:2194:G:H5'	35:BA:2195:C:OP2	2.13	0.48
26:B0:43:THR:N	35:BA:2331:G:H4'	2.29	0.48
35:BA:2359:C:H6	35:BA:2359:C:O5'	1.97	0.48
35:BA:2472:G:C5'	35:BA:2473:U:H5''	2.42	0.48
35:BA:2550:G:H2'	35:BA:2551:C:H6	1.79	0.48
35:BA:2711:A:OP1	35:BA:2712(A):A:OP1	2.31	0.48
35:BA:2791:C:H6	35:BA:2793:G:O6	1.97	0.48
35:BA:576:U:H2'	35:BA:577:G:C8	2.48	0.48
35:BA:677:A:N1	35:BA:802:A:C5	2.81	0.48
35:BA:80:G:O2'	35:BA:81:G:H5'	2.13	0.48
37:BC:67:GLY:O	37:BC:69:GLY:N	2.46	0.48
38:BD:25:THR:HG23	38:BD:25:THR:O	2.12	0.48
39:BE:3:GLY:O	39:BE:4:ILE:CB	2.62	0.48
41:BG:138:GLN:C	41:BG:140:ILE:H	2.16	0.48
41:BG:77:ILE:CD1	41:BG:82:LEU:O	2.62	0.48
42:BH:13:LYS:C	42:BH:15:VAL:N	2.65	0.48
43:BI:20:ASP:O	43:BI:21:VAL:HG13	2.14	0.48
43:BI:45:LYS:O	43:BI:48:GLU:HB3	2.13	0.48
45:BO:119:PRO:HB2	50:BT:68:TYR:HE1	1.75	0.48
46:BP:34:GLY:O	46:BP:35:HIS:C	2.51	0.48
35:BA:958:U:OP2	47:BQ:74:TYR:HE1	1.97	0.48
47:BQ:77:LYS:C	47:BQ:79:LEU:H	2.17	0.48
47:BQ:73:PRO:HA	47:BQ:93:TYR:CD2	2.48	0.48
35:BA:1652:A:H62	48:BR:11:ASN:HD21	1.61	0.48
51:BU:88:ILE:O	51:BU:90:VAL:N	2.47	0.48
35:BA:143:G:O4'	54:BX:38:GLU:HG3	2.14	0.48
56:BZ:101:PRO:O	56:BZ:102:LEU:HD23	2.13	0.48
56:BZ:146:ILE:O	56:BZ:147:GLY:C	2.51	0.48
56:BZ:91:LEU:H	56:BZ:91:LEU:HD12	1.77	0.48
2:CB:135:GLN:O	2:CB:139:LYS:HB2	2.14	0.48
2:CB:77:ALA:O	2:CB:80:ILE:HG23	2.12	0.48
2:CB:85:ALA:HB1	2:CB:92:TYR:HB3	1.96	0.48
3:CC:40:ARG:HH11	3:CC:40:ARG:HG3	1.77	0.48
4:CD:202:LEU:C	4:CD:204:ILE:N	2.66	0.48
6:CF:21:LEU:C	6:CF:25:ILE:HG12	2.34	0.48
7:CG:22:LEU:HD12	7:CG:22:LEU:O	2.13	0.48
9:CI:87:GLN:C	9:CI:89:ASN:H	2.17	0.48
11:CK:86:GLY:N	11:CK:112:THR:OG1	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:49:THR:HG22	13:CM:50:GLU:N	2.29	0.48
15:CO:74:ASP:OD2	15:CO:77:ARG:HG2	2.14	0.48
16:CP:25:ARG:HG3	16:CP:25:ARG:HH11	1.79	0.48
17:CQ:47:PRO:HG2	17:CQ:48:GLU:CD	2.34	0.48
17:CQ:70:ARG:C	17:CQ:71:PHE:HD2	2.16	0.48
1:CA:1235:U:H5'	21:CU:14:TRP:CZ2	2.48	0.48
23:CW:39:A:C2'	23:CW:40:C:H5'	2.44	0.48
23:CW:43:G:C2'	23:CW:44:A:H5''	2.44	0.48
23:CW:8:U:O2	23:CW:8:U:H2'	2.14	0.48
27:D1:71:TYR:HA	27:D1:74:VAL:CG2	2.44	0.48
29:D3:11:SER:OG	29:D3:12:PRO:HD2	2.13	0.48
29:D3:28:LEU:HA	29:D3:33:GLN:OE1	2.12	0.48
35:DA:1390:U:H6	35:DA:1390:U:OP2	1.97	0.48
35:DA:142:A:C8	35:DA:1408:C:H1'	2.49	0.48
35:DA:1465:G:H5'	35:DA:1528:A:H8	1.79	0.48
35:DA:1638:C:H2'	35:DA:1639:U:C6	2.49	0.48
35:DA:2314:C:H2'	35:DA:2315:G:C8	2.47	0.48
35:DA:2246:G:N2	35:DA:2426:A:H1'	2.29	0.48
35:DA:2650:U:H2'	35:DA:2651:C:C6	2.49	0.48
35:DA:2881:C:H2'	35:DA:2882:A:C8	2.43	0.48
35:DA:2892:A:N6	35:DA:2893:G:H21	2.12	0.48
35:DA:360:G:H2'	35:DA:361:G:C8	2.49	0.48
35:DA:365:C:H5'	35:DA:365:C:H6	1.78	0.48
35:DA:67:U:H2'	35:DA:68:G:C8	2.48	0.48
35:DA:795:C:H2'	35:DA:796:C:H6	1.78	0.48
35:DA:78:A:O2'	35:DA:79:G:H5'	2.13	0.48
36:DB:40:U:C4	36:DB:43:C:H5''	2.47	0.48
36:DB:4:C:H2'	36:DB:5:C:O4'	2.13	0.48
38:DD:96:HIS:CE1	38:DD:102:LYS:HD2	2.49	0.48
35:DA:1256:G:O2'	40:DF:82:ILE:HD12	2.14	0.48
41:DG:171:ALA:O	41:DG:172:LEU:C	2.52	0.48
41:DG:28:VAL:O	41:DG:31:VAL:HG11	2.13	0.48
41:DG:60:LEU:CA	41:DG:63:ILE:HG13	2.42	0.48
41:DG:85:GLY:O	41:DG:86:MET:HB2	2.13	0.48
41:DG:92:VAL:CG2	41:DG:93:THR:N	2.77	0.48
42:DH:87:LEU:HD23	42:DH:164:TYR:HA	1.94	0.48
42:DH:66:GLY:O	42:DH:69:ARG:N	2.47	0.48
43:DI:2:LYS:O	43:DI:3:VAL:HG13	2.14	0.48
43:DI:48:GLU:O	43:DI:51:ILE:HB	2.12	0.48
44:DN:120:LEU:HD13	44:DN:121:LYS:N	2.28	0.48
40:DF:117:ARG:CZ	46:DP:5:ASP:N	2.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:68:GLN:O	46:DP:68:GLN:HG3	2.14	0.48
48:DR:63:ARG:C	48:DR:67:LEU:HD23	2.33	0.48
50:DT:100:TYR:O	50:DT:103:ARG:N	2.47	0.48
50:DT:106:SER:O	50:DT:107:ASP:CB	2.62	0.48
35:DA:998:C:OP2	51:DU:93:LYS:HE2	2.13	0.48
51:DU:95:LEU:HA	51:DU:97:ASP:OD1	2.13	0.48
56:DZ:140:ASP:OD2	56:DZ:140:ASP:N	2.47	0.48
56:DZ:76:LEU:N	56:DZ:76:LEU:HD23	2.26	0.48
1:AA:1133:G:H2'	1:AA:1134:G:C8	2.49	0.48
1:AA:1383:C:O2'	1:AA:1384:C:H5'	2.13	0.48
1:AA:156:G:C2	1:AA:166:G:N1	2.81	0.48
1:AA:327:A:C4	1:AA:329:A:C8	3.02	0.48
1:AA:408:A:C6	1:AA:409:G:C5	3.02	0.48
1:AA:730:G:C5	1:AA:731:G:H1'	2.49	0.48
1:AA:746:A:H2'	1:AA:747:C:C6	2.49	0.48
1:AA:764:C:C2'	1:AA:765:G:H8	2.27	0.48
1:AA:836:G:C6	1:AA:851:G:C6	3.01	0.48
1:AA:940:C:H2'	1:AA:941:G:C8	2.48	0.48
1:AA:971:G:C4'	1:AA:972:C:H5''	2.42	0.48
3:AC:156:ARG:O	3:AC:158:GLY:N	2.47	0.48
3:AC:40:ARG:HB3	3:AC:44:GLU:OE2	2.14	0.48
4:AD:3:ARG:O	4:AD:4:TYR:C	2.52	0.48
6:AF:40:VAL:HG23	6:AF:62:TRP:O	2.14	0.48
8:AH:4:ASP:OD1	8:AH:6:ILE:HB	2.14	0.48
8:AH:27:PRO:HG3	8:AH:58:TYR:CE2	2.48	0.48
13:AM:11:ARG:O	13:AM:13:LYS:HG3	2.14	0.48
15:AO:62:GLN:O	15:AO:63:ARG:C	2.51	0.48
15:AO:77:ARG:O	15:AO:80:ALA:HB3	2.14	0.48
1:AA:625:G:H4'	16:AP:16:HIS:CD2	2.49	0.48
18:AR:66:LEU:HG	18:AR:70:ILE:HD11	1.96	0.48
21:AU:12:LYS:HG2	21:AU:22:ARG:HB2	1.96	0.48
23:AW:53:G:N2	23:AW:64:G:H1	2.11	0.48
25:AY:14:MET:HA	25:AY:132:ILE:CD1	2.44	0.48
27:B1:28:GLY:C	27:B1:30:VAL:N	2.52	0.48
34:B8:10:ALA:C	34:B8:12:LYS:H	2.17	0.48
34:B8:25:MET:HB2	46:BP:62:LEU:CD1	2.42	0.48
35:BA:1361:G:O2'	35:BA:1362:C:H5'	2.13	0.48
35:BA:2253:G:C6	35:BA:2254:C:N3	2.82	0.48
35:BA:2464:C:O2'	35:BA:2465:C:P	2.72	0.48
35:BA:2475:C:N4	35:BA:2529:G:H22	1.96	0.48
35:BA:2718:G:O2'	35:BA:2719:G:H5'	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:271(H):G:HO2'	35:BA:271(I):G:H8	1.62	0.48
33:B7:11:LYS:HE2	35:BA:686:G:H5''	1.96	0.48
35:BA:745:G:C2'	35:BA:746:A:H5'	2.43	0.48
35:BA:860:U:O4'	35:BA:860:U:O2	2.30	0.48
36:BB:64:C:H2'	36:BB:65:C:H6	1.77	0.48
36:BB:79:C:C2'	36:BB:80:U:H5'	2.43	0.48
37:BC:64:LEU:CD1	37:BC:66:HIS:HB2	2.43	0.48
38:BD:117:VAL:HG21	38:BD:128:GLY:C	2.34	0.48
38:BD:209:ALA:O	38:BD:210:GLY:O	2.31	0.48
39:BE:117:MET:HE3	39:BE:124:GLY:HA3	1.95	0.48
39:BE:19:ARG:O	39:BE:19:ARG:HG3	2.13	0.48
39:BE:44:TYR:CD1	39:BE:44:TYR:N	2.82	0.48
40:BF:12:LEU:O	40:BF:14:PRO:HD3	2.14	0.48
41:BG:43:LEU:CD2	41:BG:43:LEU:N	2.77	0.48
41:BG:86:MET:N	41:BG:87:PRO:CD	2.77	0.48
42:BH:153:LYS:O	42:BH:161:GLY:HA3	2.14	0.48
43:BI:54:GLN:OE1	43:BI:54:GLN:C	2.52	0.48
45:BO:98:VAL:HG22	45:BO:99:PHE:N	2.29	0.48
47:BQ:134:ARG:C	47:BQ:136:ALA:H	2.17	0.48
47:BQ:48:GLU:O	47:BQ:52:VAL:HG12	2.13	0.48
47:BQ:58:PHE:O	47:BQ:59:ARG:C	2.52	0.48
48:BR:26:LYS:HE2	48:BR:71:GLN:N	2.28	0.48
51:BU:98:LEU:C	51:BU:100:VAL:N	2.67	0.48
51:BU:110:VAL:O	51:BU:113:ALA:HB3	2.14	0.48
51:BU:29:SER:O	51:BU:30:LYS:NZ	2.41	0.48
35:BA:444:C:O5'	51:BU:2:PRO:HD3	2.12	0.48
51:BU:7:GLY:O	51:BU:8:VAL:CG2	2.62	0.48
51:BU:87:GLY:O	51:BU:88:ILE:HG23	2.14	0.48
52:BV:71:LEU:CD1	52:BV:72:VAL:H	2.18	0.48
56:BZ:6:LYS:HG3	56:BZ:62:PRO:HG3	1.96	0.48
56:BZ:79:ARG:O	56:BZ:80:ARG:HB2	2.14	0.48
1:CA:1015:A:H2'	1:CA:1016:A:C8	2.48	0.48
1:CA:112:G:C2'	1:CA:113:G:H5'	2.44	0.48
1:CA:1487:G:O2'	1:CA:1488:G:H5'	2.13	0.48
1:CA:359:U:H2'	1:CA:360:A:H8	1.79	0.48
1:CA:372:C:H4'	1:CA:373:A:OP1	2.13	0.48
1:CA:833:U:H2'	1:CA:834:C:C5	2.49	0.48
2:CB:175:ARG:O	2:CB:177:ALA:N	2.46	0.48
2:CB:19:HIS:O	2:CB:20:GLU:O	2.32	0.48
2:CB:36:ARG:O	2:CB:37:ASN:C	2.52	0.48
2:CB:39:ILE:HG22	2:CB:40:HIS:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:148:ASN:HD22	7:CG:148:ASN:N	2.11	0.48
9:CI:85:LEU:HD12	9:CI:85:LEU:C	2.33	0.48
13:CM:37:THR:O	13:CM:39:ILE:HG13	2.13	0.48
14:CN:36:PHE:O	14:CN:36:PHE:CD1	2.66	0.48
16:CP:82:GLN:O	16:CP:84:ALA:N	2.47	0.48
13:CM:84:ILE:HB	19:CS:66:MET:HE1	1.96	0.48
20:CT:22:ARG:O	20:CT:23:ARG:C	2.51	0.48
23:CW:43:G:H2'	23:CW:44:A:C4'	2.44	0.48
25:CY:10:THR:C	25:CY:12:SER:H	2.17	0.48
27:D1:88:LYS:O	27:D1:92:LYS:HB2	2.13	0.48
27:D1:93:GLU:N	27:D1:93:GLU:CD	2.67	0.48
28:D2:29:LYS:HA	28:D2:32:LEU:CD2	2.37	0.48
29:D3:48:GLU:O	29:D3:51:ALA:HB2	2.14	0.48
31:D5:31:VAL:CB	31:D5:32:PRO:HD2	2.34	0.48
35:DA:1203:G:O6	35:DA:1204:A:N6	2.47	0.48
35:DA:1324:G:H3'	35:DA:1325:G:C4'	2.43	0.48
35:DA:1515:G:O2'	35:DA:1516:C:H5'	2.14	0.48
35:DA:1812:A:C1'	38:DD:46:GLN:HE22	2.27	0.48
35:DA:2208:A:H1'	35:DA:2219:G:C5	2.49	0.48
35:DA:2681:C:C4	35:DA:2724:C:C5	3.02	0.48
35:DA:2730:C:H2'	35:DA:2731:G:H8	1.78	0.48
35:DA:431:U:H6	35:DA:431:U:O5'	1.96	0.48
35:DA:669:G:N2	35:DA:670:A:C4	2.82	0.48
35:DA:782:A:P	35:DA:782:A:H8	2.36	0.48
35:DA:830:G:N3	35:DA:2448:A:N6	2.62	0.48
35:DA:862:G:H3'	35:DA:863:A:H8	1.79	0.48
35:DA:862:G:H2'	35:DA:863:A:O4'	2.12	0.48
35:DA:90:U:O2'	35:DA:92:A:H5''	2.13	0.48
35:DA:957:A:H5'	47:DQ:76:LYS:HG3	1.96	0.48
39:DE:105:THR:HB	39:DE:197:ILE:HG12	1.96	0.48
39:DE:120:TRP:O	39:DE:122:PHE:N	2.46	0.48
40:DF:9:ILE:HG12	40:DF:14:PRO:CA	2.43	0.48
42:DH:153:LYS:O	42:DH:161:GLY:HA3	2.13	0.48
44:DN:36:GLY:HA3	44:DN:48:MET:CE	2.43	0.48
45:DO:13:ASN:HD21	45:DO:96:THR:N	2.11	0.48
45:DO:85:VAL:HG12	45:DO:86:ILE:N	2.28	0.48
47:DQ:20:ALA:CB	47:DQ:99:PRO:HG2	2.36	0.48
49:DS:87:PHE:CZ	49:DS:97:ARG:NH2	2.82	0.48
50:DT:100:TYR:CD1	50:DT:100:TYR:N	2.81	0.48
52:DV:4:ILE:CG1	52:DV:40:LEU:HD11	2.43	0.48
53:DW:20:VAL:O	53:DW:23:LEU:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:DW:65:LEU:HD23	53:DW:65:LEU:C	2.34	0.48
54:DX:58:HIS:C	54:DX:59:VAL:HG22	2.34	0.48
54:DX:72:LYS:HE3	54:DX:74:PRO:CB	2.28	0.48
56:DZ:137:ILE:HD12	56:DZ:157:LEU:HA	1.95	0.48
56:DZ:125:LEU:O	56:DZ:165:VAL:HG23	2.14	0.48
1:AA:378:G:P	16:AP:3:LYS:HZ1	2.37	0.48
1:AA:441:A:C6	1:AA:494:U:C2	3.02	0.48
2:AB:191:ASP:OD1	2:AB:192:SER:N	2.47	0.48
4:AD:196:LEU:HB3	4:AD:197:PRO:HD2	1.95	0.48
4:AD:64:LEU:HG	4:AD:65:ARG:N	2.29	0.48
4:AD:59:ARG:NH2	4:AD:66:ARG:HH22	2.04	0.48
4:AD:96:LEU:O	4:AD:98:GLU:N	2.47	0.48
5:AE:126:ARG:O	5:AE:128:PRO:N	2.47	0.48
6:AF:3:ARG:NH1	6:AF:38:GLU:OE2	2.47	0.48
7:AG:100:ALA:O	7:AG:101:LEU:C	2.51	0.48
7:AG:112:PRO:HD2	7:AG:113:GLU:OE2	2.13	0.48
11:AK:21:ILE:CD1	11:AK:82:VAL:HG13	2.36	0.48
13:AM:28:ALA:C	13:AM:30:ALA:N	2.67	0.48
16:AP:32:TYR:CE2	16:AP:35:LYS:HB2	2.49	0.48
19:AS:16:LEU:C	19:AS:20:LEU:HG	2.32	0.48
19:AS:67:VAL:HG12	19:AS:68:GLY:N	2.28	0.48
20:AT:83:ARG:CA	20:AT:86:ARG:HB3	2.42	0.48
21:AU:5:ASP:OD1	21:AU:6:ARG:N	2.47	0.48
23:AW:17:C:H5'	23:AW:18:U:C6	2.48	0.48
25:AY:108:GLU:HA	25:AY:111:ARG:HG2	1.95	0.48
25:AY:149:LEU:HB3	25:AY:153:GLU:HB2	1.95	0.48
25:AY:16:LYS:HA	25:AY:19:GLU:CG	2.43	0.48
27:B1:10:LYS:HB2	27:B1:14:VAL:N	2.29	0.48
28:B2:46:GLN:HE21	28:B2:47:ASN:CA	2.26	0.48
32:B6:28:ARG:O	32:B6:29:ASN:C	2.52	0.48
32:B6:47:THR:HG22	32:B6:49:HIS:H	1.79	0.48
35:BA:1015:G:N3	35:BA:1015:G:H2'	2.28	0.48
35:BA:1686:C:H3'	35:BA:1687:G:H8	1.78	0.48
35:BA:1686:C:H42	35:BA:1702:G:H1	1.60	0.48
35:BA:1803:A:H4'	38:BD:259:THR:CG2	2.44	0.48
35:BA:2511:U:O3'	39:BE:123:ALA:HB3	2.14	0.48
35:BA:2740:A:C2'	35:BA:2741:A:C8	2.87	0.48
35:BA:2747:G:C2	35:BA:2754:U:C4	3.02	0.48
35:BA:778:G:C5	35:BA:779:U:C4	3.02	0.48
37:BC:35:ALA:O	37:BC:36:LYS:HE3	2.14	0.48
37:BC:50:ASP:O	37:BC:52:ARG:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BC:75:LEU:HA	37:BC:94:VAL:HG22	1.95	0.48
38:BD:135:PHE:N	38:BD:135:PHE:CD1	2.72	0.48
38:BD:173:VAL:CG1	38:BD:185:VAL:O	2.62	0.48
38:BD:186:HIS:CD2	38:BD:188:GLU:HG2	2.49	0.48
39:BE:52:LEU:HB3	39:BE:76:ARG:H	1.77	0.48
40:BF:125:LEU:HD11	40:BF:199:TRP:CG	2.48	0.48
40:BF:126:VAL:O	40:BF:127:GLU:HB2	2.13	0.48
41:BG:170:ARG:HH22	41:BG:182:LYS:HZ1	1.61	0.48
41:BG:29:TRP:C	41:BG:31:VAL:N	2.66	0.48
41:BG:77:ILE:CD1	41:BG:81:LYS:O	2.62	0.48
41:BG:3:LEU:O	41:BG:8:LYS:HE2	2.14	0.48
42:BH:102:ALA:HB2	42:BH:116:GLU:HA	1.96	0.48
42:BH:43:VAL:CG1	42:BH:52:VAL:HA	2.29	0.48
44:BN:67:LEU:HD22	44:BN:88:GLU:HG2	1.96	0.48
44:BN:70:LYS:HG3	44:BN:72:TYR:HE1	1.78	0.48
44:BN:77:GLY:O	44:BN:78:TYR:CB	2.62	0.48
45:BO:61:VAL:CG1	45:BO:85:VAL:HB	2.43	0.48
46:BP:126:VAL:HG22	46:BP:145:PRO:HB3	1.96	0.48
46:BP:47:ASP:HB2	46:BP:51:PHE:HD2	1.78	0.48
49:BS:83:LYS:HA	49:BS:104:GLY:CA	2.40	0.48
50:BT:53:ARG:NH1	50:BT:53:ARG:CG	2.75	0.48
50:BT:62:THR:CG2	50:BT:75:ILE:HG13	2.42	0.48
51:BU:113:ALA:O	51:BU:116:ALA:HB3	2.14	0.48
53:BW:92:ARG:HG2	53:BW:92:ARG:NH1	2.29	0.48
54:BX:31:HIS:HD2	54:BX:33:LYS:O	1.97	0.48
54:BX:33:LYS:O	54:BX:34:ALA:C	2.52	0.48
56:BZ:139:VAL:O	56:BZ:141:VAL:N	2.45	0.48
56:BZ:149:SER:CB	56:BZ:172:ALA:O	2.62	0.48
1:CA:961:U:OP2	1:CA:1223:C:H4'	2.14	0.48
1:CA:1347:G:O2'	1:CA:1348:U:OP2	2.28	0.48
1:CA:1419:G:N2	1:CA:1482:G:H1'	2.29	0.48
1:CA:453:A:H2'	1:CA:454:C:C6	2.49	0.48
1:CA:719:C:H2'	1:CA:720:C:H5'	1.96	0.48
1:CA:788:U:C5	1:CA:789:U:C5	3.02	0.48
2:CB:111:ARG:O	2:CB:115:LEU:N	2.47	0.48
2:CB:187:LEU:HD23	2:CB:201:ILE:O	2.14	0.48
3:CC:105:GLU:HG2	3:CC:106:VAL:N	2.21	0.48
3:CC:116:VAL:O	3:CC:119:ARG:HB3	2.14	0.48
4:CD:121:VAL:HA	4:CD:126:ILE:HD12	1.96	0.48
5:CE:93:PRO:HA	5:CE:118:ILE:HD12	1.96	0.48
6:CF:12:PRO:HG3	6:CF:57:GLN:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CF:19:LEU:HD23	6:CF:19:LEU:C	2.34	0.48
8:CH:6:ILE:HD12	8:CH:6:ILE:N	2.28	0.48
9:CI:45:ALA:O	9:CI:49:PRO:HD2	2.14	0.48
11:CK:57:THR:HG23	11:CK:58:PRO:HD2	1.96	0.48
12:CL:119:LYS:H	12:CL:119:LYS:HD2	1.78	0.48
15:CO:37:ASN:N	15:CO:37:ASN:HD22	2.12	0.48
20:CT:97:ALA:O	20:CT:99:LEU:N	2.39	0.48
25:CY:144:ALA:O	25:CY:147:LEU:O	2.32	0.48
25:CY:166:ASP:O	25:CY:169:ILE:N	2.47	0.48
25:CY:171:LYS:CD	25:CY:175:LEU:HD13	2.41	0.48
29:D3:36:VAL:HG23	29:D3:36:VAL:O	2.12	0.48
34:D8:10:ALA:C	34:D8:12:LYS:H	2.17	0.48
35:DA:1456:G:C4	35:DA:1457:A:N7	2.82	0.48
35:DA:1467:C:H2'	35:DA:1468:C:H6	1.79	0.48
35:DA:1479:G:H2'	35:DA:1480:G:O4'	2.13	0.48
35:DA:2472:G:C5'	35:DA:2473:U:H5''	2.41	0.48
35:DA:2759:G:H5'	35:DA:2759:G:H8	1.79	0.48
35:DA:2845:G:H5''	50:DT:55:ASN:HA	1.96	0.48
35:DA:2846:G:H2'	35:DA:2847:U:H6	1.79	0.48
35:DA:614:U:H4'	35:DA:614(C):A:H62	1.79	0.48
35:DA:658:C:H2'	35:DA:659:C:C6	2.49	0.48
35:DA:692:C:N3	35:DA:771:G:C2	2.81	0.48
35:DA:778:G:C6	35:DA:779:U:N3	2.81	0.48
35:DA:810:U:C2	46:DP:31:ALA:O	2.66	0.48
35:DA:958:U:OP2	47:DQ:74:TYR:HE1	1.97	0.48
36:DB:5:C:H2'	36:DB:6:C:H6	1.79	0.48
37:DC:44:HIS:HD2	37:DC:175:VAL:CA	2.27	0.48
38:DD:172:TYR:CD1	38:DD:186:HIS:CA	2.95	0.48
39:DE:173:VAL:O	39:DE:174:ASP:C	2.52	0.48
40:DF:6:VAL:CG2	40:DF:124:LEU:HA	2.41	0.48
40:DF:186:ILE:HG23	40:DF:192:LEU:HD12	1.95	0.48
40:DF:202:PHE:CE1	40:DF:206:ILE:HD13	2.45	0.48
44:DN:64:GLY:O	44:DN:65:LYS:HE3	2.13	0.48
45:DO:112:MET:O	45:DO:113:LYS:C	2.51	0.48
45:DO:2:ILE:CD1	45:DO:6:THR:HG21	2.43	0.48
46:DP:106:LEU:CD1	46:DP:112:LEU:HB2	2.42	0.48
46:DP:62:LEU:CD1	46:DP:62:LEU:N	2.64	0.48
48:DR:118:GLU:HA	48:DR:118:GLU:OE1	2.12	0.48
49:DS:92:TYR:C	49:DS:92:TYR:CD1	2.87	0.48
50:DT:52:ILE:CG2	50:DT:61:PHE:HB2	2.42	0.48
51:DU:98:LEU:O	51:DU:101:ARG:O	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:DU:33:ARG:O	51:DU:36:ARG:N	2.46	0.48
51:DU:59:ARG:C	51:DU:61:TRP:N	2.66	0.48
51:DU:78:THR:C	51:DU:80:ILE:N	2.66	0.48
52:DV:75:PHE:HB2	52:DV:87:HIS:CB	2.43	0.48
53:DW:64:MET:CE	53:DW:109:GLU:HG3	2.44	0.48
35:DA:1615:C:O2	53:DW:87:PRO:HG2	2.14	0.48
53:DW:96:ILE:HG13	53:DW:97:LYS:H	1.79	0.48
35:DA:143:G:C4'	54:DX:38:GLU:HG3	2.44	0.48
54:DX:83:VAL:O	54:DX:83:VAL:HG23	2.13	0.48
56:DZ:150:LEU:HD23	56:DZ:171:ILE:CG1	2.44	0.48
1:AA:124:G:C6	1:AA:125:U:C4	3.01	0.48
1:AA:1316:G:H2'	1:AA:1317:C:C5'	2.43	0.48
1:AA:542:G:H5'	4:AD:41:GLY:HA3	1.96	0.48
1:AA:673:G:H2'	1:AA:674:G:H8	1.71	0.48
1:AA:956:U:H2'	1:AA:957:U:H6	1.79	0.48
2:AB:92:TYR:CE2	2:AB:151:GLY:HA3	2.49	0.48
4:AD:131:ARG:HD3	4:AD:131:ARG:N	2.29	0.48
6:AF:55:ASP:OD2	6:AF:86:ARG:NH2	2.47	0.48
7:AG:122:HIS:O	7:AG:125:MET:N	2.47	0.48
8:AH:10:LEU:CD2	8:AH:10:LEU:H	2.17	0.48
8:AH:68:ARG:HG3	8:AH:69:ARG:N	2.24	0.48
11:AK:48:ILE:HD13	11:AK:48:ILE:N	2.29	0.48
16:AP:25:ARG:HH11	16:AP:25:ARG:HG3	1.78	0.48
1:AA:1235:U:H5'	21:AU:14:TRP:CZ2	2.49	0.48
21:AU:20:LYS:HD3	21:AU:21:TYR:HE1	1.77	0.48
7:AG:143:ARG:CZ	23:AW:43:G:H5'	2.44	0.48
31:B5:20:ARG:HH12	53:BW:15:ARG:HH21	1.55	0.48
35:BA:1132:A:H2'	35:BA:1133:U:C6	2.48	0.48
35:BA:1198:U:H2'	35:BA:1199:U:H6	1.79	0.48
35:BA:1203:G:O6	35:BA:1204:A:N6	2.46	0.48
35:BA:1233:C:H2'	35:BA:1234:U:H5'	1.95	0.48
35:BA:1374:G:H2'	35:BA:1375:C:H6	1.77	0.48
35:BA:1496:A:H8	35:BA:1577:C:HO2'	1.62	0.48
35:BA:1578:U:OP2	35:BA:1578:U:H6	1.96	0.48
35:BA:1688:U:O2	35:BA:1700:A:H8	1.97	0.48
35:BA:2124:G:C2'	35:BA:2125:G:H5'	2.44	0.48
35:BA:2291:U:H2'	35:BA:2292:C:C6	2.49	0.48
35:BA:2312:U:H2'	35:BA:2313:C:C5'	2.43	0.48
35:BA:2729:G:H2'	35:BA:2730:C:C6	2.49	0.48
35:BA:514:A:C2	35:BA:515:A:C4	3.02	0.48
35:BA:684:G:N2	35:BA:787:U:H2'	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:117:VAL:CG2	38:BD:118:VAL:H	2.26	0.48
38:BD:130:ALA:HB2	38:BD:192:THR:HA	1.95	0.48
38:BD:199:ALA:O	38:BD:201:HIS:N	2.47	0.48
39:BE:13:ARG:HA	39:BE:22:PRO:HA	1.96	0.48
40:BF:9:ILE:HG23	40:BF:14:PRO:HA	1.96	0.48
41:BG:106:LEU:HD12	41:BG:110:ALA:HB3	1.94	0.48
42:BH:105:LEU:N	42:BH:105:LEU:CD2	2.74	0.48
43:BI:119:PRO:O	43:BI:120:ILE:HB	2.14	0.48
47:BQ:101:ARG:HG3	47:BQ:101:ARG:NH1	2.29	0.48
47:BQ:119:ARG:HG2	47:BQ:120:ILE:CD1	2.41	0.48
48:BR:88:ARG:NH2	48:BR:89:ASP:OD1	2.47	0.48
49:BS:24:LEU:O	49:BS:86:ALA:HB3	2.13	0.48
50:BT:10:VAL:O	50:BT:11:GLU:C	2.52	0.48
50:BT:121:ILE:CG2	50:BT:122:ASP:N	2.77	0.48
52:BV:66:ARG:HB2	52:BV:95:LEU:H	1.78	0.48
55:BY:45:VAL:HG22	55:BY:62:GLU:CB	2.36	0.48
56:BZ:126:VAL:HG12	56:BZ:162:GLU:O	2.14	0.48
1:CA:1164:G:H2'	1:CA:1165:C:H5'	1.96	0.48
1:CA:1293:G:O2'	1:CA:1294:G:H8	1.96	0.48
1:CA:1237:C:H4'	1:CA:1334:G:N2	2.29	0.48
1:CA:658:G:O2'	1:CA:659:U:H5'	2.14	0.48
1:CA:689:C:H2'	1:CA:689:C:O2	2.13	0.48
1:CA:782:A:N6	1:CA:801:U:C6	2.81	0.48
1:CA:81:U:H2'	1:CA:82:U:C6	2.49	0.48
1:CA:978:A:O2'	1:CA:1322:C:N3	2.47	0.48
2:CB:233:SER:C	2:CB:235:SER:H	2.18	0.48
2:CB:42:ILE:CD1	2:CB:203:GLY:H	2.27	0.48
4:CD:192:GLU:O	4:CD:194:LEU:N	2.47	0.48
5:CE:51:VAL:HB	5:CE:52:PRO:CD	2.37	0.48
9:CI:119:ALA:O	9:CI:120:ARG:HB2	2.14	0.48
9:CI:4:TYR:CE1	9:CI:21:PRO:HD3	2.48	0.48
11:CK:58:PRO:HA	11:CK:90:GLY:CA	2.43	0.48
15:CO:17:ARG:HG3	15:CO:17:ARG:NH1	2.24	0.48
17:CQ:99:SER:C	17:CQ:100:LYS:HE2	2.35	0.48
17:CQ:82:MET:HA	17:CQ:85:VAL:CG2	2.44	0.48
18:CR:34:TYR:O	18:CR:35:ARG:HG2	2.14	0.48
19:CS:36:ARG:NH1	19:CS:36:ARG:HB3	2.29	0.48
25:CY:107:THR:O	25:CY:111:ARG:HG3	2.14	0.48
27:D1:87:PRO:CD	27:D1:88:LYS:H	2.27	0.48
28:D2:30:ARG:HH11	28:D2:30:ARG:CG	2.27	0.48
28:D2:38:GLN:C	28:D2:40:SER:N	2.67	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:D5:16:ARG:HG2	31:D5:16:ARG:NH1	2.24	0.48
31:D5:42:PRO:C	31:D5:43:HIS:HD2	2.17	0.48
35:DA:1613:G:H3'	35:DA:1614:A:H5''	1.95	0.48
35:DA:1660:C:H5'	35:DA:2712(A):A:N6	2.27	0.48
35:DA:1891:G:C2	35:DA:1892:C:C2	3.02	0.48
27:D1:37:ILE:CD1	35:DA:2080:G:OP1	2.62	0.48
35:DA:301:G:H1'	35:DA:302:C:C6	2.49	0.48
35:DA:363(F):A:O2'	35:DA:364:C:H5	1.96	0.48
35:DA:374:A:C2	35:DA:375:C:H1'	2.49	0.48
35:DA:449:A:C2'	35:DA:450:G:H5'	2.43	0.48
35:DA:467:G:H2'	35:DA:468:G:C8	2.42	0.48
35:DA:537:C:H3'	35:DA:538:G:H8	1.78	0.48
35:DA:565:C:H2'	35:DA:566:U:H6	1.79	0.48
35:DA:924:C:O2'	35:DA:925:C:H5'	2.14	0.48
35:DA:968:G:H2'	35:DA:969:U:C6	2.48	0.48
35:DA:970:C:H2'	35:DA:971:C:C6	2.48	0.48
35:DA:998:C:OP2	51:DU:93:LYS:NZ	2.47	0.48
36:DB:66:A:C2	36:DB:109:C:C2	3.02	0.48
37:DC:51:PRO:O	37:DC:52:ARG:HB2	2.14	0.48
35:DA:1490:A:H2	38:DD:75:ILE:HD12	1.79	0.48
39:DE:24:THR:HB	39:DE:186:GLY:HA2	1.95	0.48
40:DF:17:ARG:HH11	40:DF:17:ARG:HG3	1.78	0.48
40:DF:52:LYS:O	40:DF:88:VAL:HG12	2.14	0.48
42:DH:78:GLY:O	42:DH:136:ILE:HG23	2.14	0.48
44:DN:26:LEU:HG	44:DN:27:ALA:N	2.27	0.48
46:DP:41:ARG:CD	46:DP:41:ARG:N	2.77	0.48
47:DQ:26:TYR:O	47:DQ:27:VAL:C	2.51	0.48
47:DQ:28:ALA:HB3	47:DQ:105:GLU:CD	2.34	0.48
35:DA:2880:C:H1'	48:DR:92:GLY:O	2.14	0.48
49:DS:38:GLN:CD	49:DS:47:THR:HG23	2.33	0.48
50:DT:45:PHE:HE2	50:DT:63:VAL:HG23	1.79	0.48
51:DU:113:ALA:O	51:DU:116:ALA:HB3	2.13	0.48
52:DV:36:PRO:HD2	52:DV:60:GLU:O	2.13	0.48
53:DW:31:GLU:O	53:DW:32:ALA:C	2.53	0.48
55:DY:18:GLY:C	55:DY:20:TYR:N	2.67	0.48
1:AA:636:U:H2'	1:AA:637:G:H8	1.72	0.47
1:AA:959:A:C2'	1:AA:960:U:H4'	2.44	0.47
2:AB:81:VAL:HG22	2:AB:215:LEU:CG	2.44	0.47
4:AD:101:LEU:CD2	4:AD:121:VAL:HG13	2.44	0.47
4:AD:68:TYR:CE2	4:AD:97:LEU:HD22	2.48	0.47
7:AG:27:ILE:HD12	7:AG:27:ILE:H	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:110:ALA:HA	8:AH:136:GLU:HA	1.96	0.47
10:AJ:39:PRO:HB3	10:AJ:70:ARG:CZ	2.44	0.47
14:AN:4:LYS:HA	14:AN:7:ILE:HD11	1.96	0.47
15:AO:3:ILE:O	15:AO:3:ILE:HG13	2.13	0.47
16:AP:23:ASP:OD2	16:AP:25:ARG:NH2	2.47	0.47
22:AV:29:G:N1	22:AV:30:A:C5	2.82	0.47
22:AV:29:G:C2	22:AV:30:A:C8	3.02	0.47
23:AW:23:G:H2'	23:AW:24:C:C5'	2.44	0.47
23:AW:8:U:H3'	23:AW:8:U:OP2	2.14	0.47
25:AY:10:THR:HG21	25:AY:164:ILE:HG21	1.96	0.47
25:AY:65:THR:CG2	25:AY:66:LEU:N	2.77	0.47
27:B1:9:GLY:O	27:B1:10:LYS:HB3	2.14	0.47
28:B2:14:ARG:HE	28:B2:14:ARG:H	1.59	0.47
29:B3:32:GLN:HA	29:B3:32:GLN:OE1	2.13	0.47
32:B6:11:LEU:O	32:B6:24:GLU:N	2.47	0.47
35:BA:1202:C:C2'	35:BA:1203:G:H5'	2.44	0.47
35:BA:1468:C:O2'	35:BA:1469:A:H5'	2.14	0.47
35:BA:1550:C:O2'	35:BA:1551:C:H5'	2.14	0.47
35:BA:1831:G:O2'	35:BA:1832:C:H5'	2.14	0.47
35:BA:184:C:O2'	35:BA:185:U:H5'	2.13	0.47
35:BA:197:A:C8	35:BA:197:A:H5'	2.49	0.47
35:BA:238:C:H2'	35:BA:239:U:H6	1.79	0.47
35:BA:412:A:N6	35:BA:2411:A:H2'	2.29	0.47
35:BA:2581:G:N2	35:BA:2581:G:OP2	2.45	0.47
35:BA:750:A:H2'	35:BA:751:A:H5''	1.95	0.47
35:BA:779:U:OP1	38:BD:49:ILE:HG23	2.14	0.47
35:BA:826:U:C2	35:BA:828:U:H1'	2.49	0.47
39:BE:4:ILE:HD12	39:BE:31:CYS:SG	2.54	0.47
40:BF:178:PRO:CG	40:BF:179:GLU:N	2.76	0.47
40:BF:34:TRP:HA	40:BF:37:VAL:HG23	1.95	0.47
41:BG:25:TYR:CE2	41:BG:31:VAL:HA	2.49	0.47
42:BH:38:SER:C	42:BH:40:GLU:H	2.18	0.47
42:BH:60:ARG:O	42:BH:64:LEU:HG	2.14	0.47
43:BI:69:LYS:C	43:BI:71:ILE:H	2.18	0.47
43:BI:79:ILE:CB	43:BI:81:VAL:HG23	2.43	0.47
44:BN:120:LEU:HD13	44:BN:121:LYS:N	2.29	0.47
44:BN:64:GLY:O	44:BN:65:LYS:HE3	2.13	0.47
45:BO:59:LYS:O	45:BO:86:ILE:HG23	2.14	0.47
49:BS:103:GLU:O	49:BS:105:ALA:N	2.46	0.47
49:BS:95:HIS:O	49:BS:98:VAL:HG23	2.14	0.47
50:BT:102:ILE:HB	50:BT:110:ILE:HD11	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:108:ARG:HH11	50:BT:108:ARG:HB3	1.79	0.47
50:BT:78:LEU:O	50:BT:78:LEU:HD23	2.14	0.47
51:BU:86:ALA:HB2	51:BU:116:ALA:CB	2.43	0.47
35:BA:58:G:OP1	54:BX:72:LYS:HA	2.14	0.47
55:BY:68:HIS:HB3	55:BY:71:LYS:HZ3	1.79	0.47
56:BZ:23:LYS:O	56:BZ:41:LEU:HD21	2.14	0.47
1:CA:1146:A:H2'	1:CA:1147:C:O5'	2.13	0.47
1:CA:1273:G:H2'	1:CA:1274:G:O4'	2.14	0.47
1:CA:1396:A:C4'	1:CA:1398:A:H1'	2.44	0.47
1:CA:1422:G:H2'	1:CA:1423:G:H8	1.79	0.47
1:CA:1519:A:H3'	1:CA:1520:G:O4'	2.13	0.47
1:CA:767:A:H2'	1:CA:768:A:H8	1.79	0.47
2:CB:19:HIS:CA	2:CB:39:ILE:HD13	2.39	0.47
3:CC:156:ARG:O	3:CC:158:GLY:N	2.47	0.47
3:CC:173:VAL:O	3:CC:173:VAL:CG1	2.61	0.47
3:CC:41:GLY:O	3:CC:45:LYS:HG3	2.15	0.47
4:CD:202:LEU:O	4:CD:205:GLU:N	2.47	0.47
4:CD:30:LYS:O	4:CD:32:ALA:N	2.47	0.47
4:CD:52:SER:OG	4:CD:55:ALA:N	2.39	0.47
5:CE:121:LYS:HD2	5:CE:122:GLU:H	1.79	0.47
7:CG:105:VAL:O	7:CG:108:ALA:HB3	2.13	0.47
7:CG:156:TRP:HD1	7:CG:156:TRP:H	1.62	0.47
7:CG:36:LYS:O	7:CG:39:ALA:HB3	2.14	0.47
12:CL:32:PHE:HE1	12:CL:86:ARG:HG3	1.79	0.47
13:CM:14:ARG:HG3	13:CM:17:VAL:CG2	2.43	0.47
16:CP:17:TYR:N	16:CP:17:TYR:CD1	2.82	0.47
19:CS:9:VAL:O	19:CS:10:PHE:C	2.52	0.47
26:D0:17:GLN:HG2	35:DA:2261:C:P	2.54	0.47
27:D1:86:SER:N	27:D1:87:PRO:CD	2.75	0.47
27:D1:85:LEU:C	27:D1:87:PRO:CD	2.81	0.47
32:D6:27:LYS:HE2	35:DA:2285:C:H5	1.78	0.47
32:D6:32:ASN:O	32:D6:33:LYS:HB2	2.14	0.47
35:DA:11:G:O2'	35:DA:12:U:H5'	2.14	0.47
35:DA:1260:G:H2'	35:DA:1261:C:H6	1.79	0.47
35:DA:1446:C:O2'	35:DA:1447:G:H5'	2.14	0.47
35:DA:1484:G:H3'	35:DA:1485:G:C5'	2.44	0.47
35:DA:120:U:H1'	35:DA:149:A:N7	2.28	0.47
35:DA:1504:C:O2'	35:DA:1505:C:C5'	2.62	0.47
35:DA:1806:C:N4	35:DA:1812:A:N6	2.62	0.47
35:DA:1817:G:O2'	35:DA:1818:U:H5'	2.13	0.47
35:DA:1902:C:H2'	35:DA:1903:G:O4'	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2253:G:C6	35:DA:2254:C:N3	2.82	0.47
35:DA:2323:G:H2'	35:DA:2324:C:O4'	2.14	0.47
35:DA:2658:C:H41	35:DA:2664:G:N2	2.12	0.47
35:DA:2694:G:O2'	35:DA:2695:C:H5'	2.14	0.47
35:DA:2708:G:H1'	48:DR:71:GLN:OE1	2.13	0.47
35:DA:271(A):A:H5''	35:DA:271(B):C:C5	2.49	0.47
35:DA:2851:A:H2'	35:DA:2852:G:C8	2.48	0.47
35:DA:297:C:H2'	35:DA:298:G:O4'	2.13	0.47
35:DA:704:G:N3	35:DA:726:G:C2	2.82	0.47
35:DA:806:C:O2'	35:DA:807:U:H5'	2.14	0.47
35:DA:892:G:N3	35:DA:892:G:H3'	2.29	0.47
37:DC:65:PRO:HG2	37:DC:189:ILE:CB	2.44	0.47
38:DD:106:ILE:HG23	38:DD:106:ILE:O	2.13	0.47
35:DA:1971:A:C4	38:DD:241:PRO:HB3	2.48	0.47
39:DE:179:GLU:O	39:DE:180:ASN:HB3	2.13	0.47
39:DE:79:ARG:NH1	39:DE:79:ARG:HG2	2.29	0.47
40:DF:125:LEU:HD11	40:DF:199:TRP:CD1	2.48	0.47
41:DG:122:PRO:O	41:DG:124:SER:N	2.47	0.47
42:DH:19:VAL:CG1	42:DH:44:VAL:HG22	2.44	0.47
42:DH:61:HIS:C	42:DH:63:SER:N	2.67	0.47
44:DN:17:ASP:OD2	44:DN:17:ASP:O	2.32	0.47
46:DP:101:VAL:HG13	46:DP:102:ARG:N	2.21	0.47
47:DQ:131:ILE:H	47:DQ:131:ILE:HD13	1.79	0.47
47:DQ:76:LYS:N	47:DQ:88:GLY:HA3	2.29	0.47
48:DR:55:ALA:HA	48:DR:80:PHE:CZ	2.48	0.47
49:DS:103:GLU:O	49:DS:105:ALA:N	2.47	0.47
51:DU:52:ARG:O	51:DU:53:ARG:C	2.53	0.47
52:DV:3:ALA:O	52:DV:13:ARG:HA	2.14	0.47
54:DX:29:TRP:HE3	54:DX:76:ARG:HB3	1.79	0.47
35:DA:58:G:OP1	54:DX:72:LYS:HA	2.13	0.47
55:DY:2:ARG:C	55:DY:4:LYS:N	2.65	0.47
56:DZ:24:LEU:HD21	56:DZ:86:VAL:CG2	2.44	0.47
1:AA:1438:G:N1	1:AA:1464:G:C2	2.82	0.47
1:AA:343:U:O2'	1:AA:344:A:H2'	2.15	0.47
1:AA:668:G:O2'	1:AA:669:U:H5'	2.14	0.47
1:AA:693:G:H2'	1:AA:694:A:O4'	2.12	0.47
1:AA:989:C:O2'	1:AA:990:C:H5'	2.13	0.47
4:AD:45:GLN:O	4:AD:46:LYS:HG3	2.14	0.47
4:AD:79:PHE:C	4:AD:79:PHE:CD2	2.87	0.47
6:AF:62:TRP:CH2	6:AF:64:GLN:HB2	2.48	0.47
7:AG:150:ALA:C	7:AG:152:ALA:N	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:69:ARG:HB2	8:AH:74:PRO:HA	1.96	0.47
11:AK:122:LYS:O	11:AK:126:ARG:HB3	2.13	0.47
13:AM:117:VAL:CG1	13:AM:118:ALA:N	2.77	0.47
13:AM:14:ARG:HG3	13:AM:17:VAL:CG2	2.44	0.47
13:AM:54:VAL:O	13:AM:58:GLU:N	2.46	0.47
18:AR:66:LEU:O	18:AR:70:ILE:HG13	2.13	0.47
20:AT:71:THR:HB	20:AT:72:LEU:H	1.55	0.47
26:B0:7:LEU:CA	47:BQ:83:MET:SD	3.02	0.47
34:B8:39:LYS:HD3	34:B8:39:LYS:C	2.34	0.47
35:BA:1022:G:N2	35:BA:1142(A):A:C2	2.82	0.47
35:BA:1184:G:C2'	35:BA:1185:C:H5'	2.44	0.47
35:BA:1613:G:N1	35:BA:1619:G:C5	2.82	0.47
35:BA:1682:G:O2'	35:BA:1683:C:H5'	2.14	0.47
35:BA:2223:G:O2'	35:BA:2224:G:H5'	2.13	0.47
35:BA:602:G:N3	35:BA:602:G:H2'	2.29	0.47
35:BA:705:A:C2	35:BA:706:A:C4	3.03	0.47
36:BB:66:A:N6	36:BB:108:U:H2'	2.25	0.47
36:BB:25:A:H2'	36:BB:26:A:O4'	2.13	0.47
36:BB:73:A:H2'	36:BB:74:U:O4'	2.14	0.47
38:BD:267:SER:HA	38:BD:270:ILE:HG13	1.95	0.47
35:BA:1816:G:N7	38:BD:62:TYR:CE1	2.82	0.47
38:BD:95:LEU:N	38:BD:95:LEU:HD12	2.29	0.47
39:BE:181:LEU:O	39:BE:182:LEU:CB	2.61	0.47
40:BF:24:LEU:CD2	40:BF:24:LEU:N	2.77	0.47
40:BF:65:TRP:CH2	40:BF:75:HIS:CD2	2.95	0.47
41:BG:7:LEU:CA	41:BG:10:LYS:HB2	2.37	0.47
41:BG:76:SER:HB2	41:BG:83:ARG:HB3	1.94	0.47
42:BH:125:VAL:HG12	42:BH:127:GLU:O	2.15	0.47
42:BH:109:PHE:HE1	42:BH:152:ARG:NE	2.12	0.47
42:BH:155:SER:OG	42:BH:156:ALA:N	2.44	0.47
42:BH:157:TYR:CD1	42:BH:170:ARG:O	2.66	0.47
43:BI:145:VAL:CG1	43:BI:146:ALA:N	2.74	0.47
43:BI:47:LEU:O	43:BI:50:ARG:N	2.41	0.47
43:BI:77:LEU:HD21	43:BI:101:LEU:HD13	1.95	0.47
45:BO:64:ARG:CB	45:BO:64:ARG:HH11	2.27	0.47
47:BQ:34:LEU:HG	47:BQ:103:MET:HB2	1.96	0.47
50:BT:110:ILE:CG2	50:BT:111:ARG:N	2.77	0.47
50:BT:30:VAL:O	50:BT:31:SER:HB3	2.14	0.47
50:BT:79:HIS:O	50:BT:80:SER:HB2	2.14	0.47
51:BU:28:ARG:HA	51:BU:34:LYS:HB3	1.96	0.47
35:BA:993:G:H1'	52:BV:91:TYR:CE1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:2:ARG:C	55:BY:4:LYS:N	2.67	0.47
55:BY:31:LEU:HD12	55:BY:34:LYS:H	1.77	0.47
56:BZ:138:GLU:OE2	56:BZ:138:GLU:N	2.47	0.47
56:BZ:20:ARG:HG3	56:BZ:20:ARG:NH1	2.29	0.47
56:BZ:61:LEU:CD1	56:BZ:65:GLN:HB3	2.43	0.47
1:CA:1016:A:H2'	1:CA:1017:G:H5'	1.96	0.47
1:CA:1120:G:H1	1:CA:1153:C:H42	1.61	0.47
1:CA:1338:G:H21	22:CV:40:C:H1'	1.80	0.47
1:CA:241:C:O2'	1:CA:242:C:H5'	2.14	0.47
1:CA:313:A:H2'	1:CA:314:C:H6	1.76	0.47
1:CA:410:G:N1	1:CA:429:U:O2	2.46	0.47
1:CA:545:C:O2'	1:CA:546:G:H5'	2.15	0.47
1:CA:55:A:N7	1:CA:56:U:C5	2.82	0.47
1:CA:872:A:C2	1:CA:874:G:C6	3.02	0.47
1:CA:880:C:H2'	1:CA:881:G:H8	1.79	0.47
2:CB:170:GLU:C	2:CB:172:ILE:N	2.68	0.47
2:CB:191:ASP:OD1	2:CB:192:SER:N	2.47	0.47
2:CB:80:ILE:HG13	2:CB:81:VAL:HG23	1.95	0.47
2:CB:82:ARG:HA	2:CB:92:TYR:CE1	2.49	0.47
3:CC:11:ARG:NH1	3:CC:11:ARG:HG2	2.29	0.47
3:CC:134:ILE:HG22	3:CC:168:ALA:CB	2.43	0.47
3:CC:84:ILE:HA	3:CC:87:LEU:HB2	1.95	0.47
6:CF:62:TRP:CH2	6:CF:64:GLN:HB2	2.49	0.47
7:CG:104:LEU:H	7:CG:104:LEU:CD2	2.27	0.47
7:CG:148:ASN:O	7:CG:150:ALA:N	2.47	0.47
8:CH:127:LEU:HD12	8:CH:129:VAL:HG13	1.97	0.47
8:CH:49:GLU:HG2	8:CH:62:TYR:HE2	1.79	0.47
9:CI:37:PHE:HB3	9:CI:43:ALA:HB2	1.96	0.47
9:CI:7:THR:H	9:CI:83:ARG:HD2	1.79	0.47
10:CJ:39:PRO:HB3	10:CJ:70:ARG:CZ	2.44	0.47
12:CL:117:ARG:O	12:CL:119:LYS:O	2.31	0.47
14:CN:39:LEU:CD1	14:CN:47:LEU:HD12	2.44	0.47
16:CP:74:LEU:O	16:CP:79:VAL:HG23	2.14	0.47
17:CQ:45:HIS:NE2	17:CQ:47:PRO:HB3	2.29	0.47
17:CQ:92:ARG:O	17:CQ:95:TYR:N	2.37	0.47
20:CT:56:MET:O	20:CT:60:GLU:HB2	2.13	0.47
26:D0:43:THR:O	26:D0:43:THR:HG23	2.13	0.47
29:D3:4:LEU:HD23	29:D3:4:LEU:C	2.34	0.47
33:D7:19:ARG:HG2	33:D7:19:ARG:NH1	2.29	0.47
35:DA:1278:A:O2'	35:DA:1279:G:H5'	2.14	0.47
35:DA:1441:G:H2'	35:DA:1442:G:H8	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1450:G:H1	35:DA:1462:C:H42	1.61	0.47
35:DA:1444:G:N2	35:DA:1548:C:C2	2.82	0.47
35:DA:1992:G:C2	35:DA:1997:G:C6	3.02	0.47
35:DA:2069:G:C2'	35:DA:2070:G:H5'	2.43	0.47
35:DA:259:G:C2	35:DA:260:G:C8	3.02	0.47
35:DA:2740:A:C6	35:DA:2764:A:C8	3.01	0.47
35:DA:283:A:O2'	35:DA:284:U:OP1	2.25	0.47
35:DA:290:G:N2	35:DA:291:C:H1'	2.29	0.47
35:DA:982:C:O5'	35:DA:982:C:H6	1.97	0.47
37:DC:64:LEU:CD1	37:DC:66:HIS:HB2	2.44	0.47
37:DC:67:GLY:O	37:DC:69:GLY:N	2.46	0.47
35:DA:1659:U:OP2	39:DE:132:HIS:CE1	2.67	0.47
40:DF:139:PHE:HB3	40:DF:166:ALA:HB1	1.95	0.47
40:DF:164:ARG:HG3	40:DF:175:THR:HG1	1.76	0.47
41:DG:130:ASN:HD22	41:DG:160:VAL:HA	1.79	0.47
41:DG:76:SER:C	41:DG:77:ILE:HG13	2.34	0.47
45:DO:26:LYS:O	45:DO:27:GLY:O	2.32	0.47
47:DQ:101:ARG:HG3	47:DQ:101:ARG:NH1	2.29	0.47
28:D2:22:GLU:HB3	54:DX:5:TYR:CD1	2.48	0.47
55:DY:60:PHE:HD2	55:DY:60:PHE:O	1.96	0.47
1:AA:1072:G:H2'	1:AA:1073:U:H6	1.78	0.47
1:AA:1438:G:C2	1:AA:1464:G:C2	3.01	0.47
1:AA:975:A:H4'	1:AA:976:G:C5'	2.28	0.47
1:AA:619:U:C2	4:AD:135:LEU:HD23	2.49	0.47
6:AF:68:PRO:HG3	6:AF:71:ARG:HE	1.79	0.47
7:AG:148:ASN:HD22	7:AG:148:ASN:N	2.12	0.47
9:AI:50:LEU:HB3	9:AI:55:ALA:CB	2.44	0.47
12:AL:76:ASN:O	12:AL:77:LEU:HD23	2.14	0.47
13:AM:114:ARG:C	13:AM:116:THR:H	2.17	0.47
15:AO:15:PHE:O	15:AO:16:ALA:O	2.31	0.47
17:AQ:10:VAL:HG11	17:AQ:53:LEU:HA	1.96	0.47
17:AQ:59:ILE:CD1	17:AQ:73:VAL:HA	2.44	0.47
18:AR:74:ARG:HG3	18:AR:74:ARG:NH1	2.29	0.47
1:AA:1242:C:P	21:AU:10:ARG:HH12	2.37	0.47
30:B4:45:GLY:C	30:B4:47:GLN:N	2.66	0.47
31:B5:40:LYS:HZ3	31:B5:50:GLY:HA2	1.79	0.47
34:B8:58:ILE:C	34:B8:61:LEU:HG	2.34	0.47
35:BA:118:A:OP2	35:BA:119:A:H5''	2.14	0.47
35:BA:2069:G:C2	35:BA:2070:G:C8	3.01	0.47
35:BA:2225:A:H4'	35:BA:2226:C:C5'	2.43	0.47
35:BA:2678:C:O2'	35:BA:2679:A:H5'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:874:G:H2'	35:BA:875:G:H8	1.78	0.47
36:BB:40:U:H2'	36:BB:41:U:OP1	2.14	0.47
37:BC:77:ILE:HG13	37:BC:96:GLY:O	2.15	0.47
38:BD:268:ARG:HB2	38:BD:268:ARG:CZ	2.44	0.47
40:BF:107:LYS:O	40:BF:108:LYS:C	2.52	0.47
41:BG:38:VAL:O	41:BG:158:ALA:CB	2.62	0.47
41:BG:76:SER:CA	41:BG:84:LYS:H	2.26	0.47
43:BI:94:ALA:CB	43:BI:114:LEU:HD12	2.44	0.47
35:BA:2415:G:C4'	46:BP:66:GLY:HA3	2.43	0.47
47:BQ:82:ARG:CG	47:BQ:82:ARG:HH11	2.22	0.47
48:BR:20:LEU:CD1	48:BR:20:LEU:C	2.75	0.47
48:BR:4:LEU:C	48:BR:6:SER:H	2.16	0.47
49:BS:25:ARG:NH2	49:BS:89:ARG:NH1	2.60	0.47
49:BS:83:LYS:O	49:BS:85:VAL:N	2.48	0.47
51:BU:52:ARG:O	51:BU:53:ARG:C	2.52	0.47
51:BU:92:ARG:CZ	52:BV:11:GLN:HG2	2.44	0.47
51:BU:98:LEU:O	51:BU:100:VAL:N	2.47	0.47
52:BV:45:THR:HG22	52:BV:45:THR:O	2.13	0.47
53:BW:106:ILE:O	53:BW:106:ILE:HG13	2.13	0.47
53:BW:37:ARG:HG3	53:BW:37:ARG:NH1	2.29	0.47
54:BX:75:ASP:O	54:BX:76:ARG:HG3	2.14	0.47
1:CA:1129:C:H4'	1:CA:1130:A:H5'	1.96	0.47
1:CA:1401:G:C2	1:CA:1402:C:H1'	2.49	0.47
1:CA:255:G:H2'	1:CA:256:U:C6	2.49	0.47
1:CA:332:G:H2'	1:CA:333:G:C8	2.46	0.47
1:CA:352:C:O2	1:CA:352:C:H2'	2.13	0.47
1:CA:35:G:H2'	1:CA:36:C:C6	2.49	0.47
1:CA:425:G:H2'	1:CA:426:G:C8	2.46	0.47
1:CA:695:A:O2'	1:CA:696:A:H5'	2.14	0.47
1:CA:746:A:H2'	1:CA:747:C:C6	2.50	0.47
1:CA:93:G:C6	1:CA:96:U:C4	3.02	0.47
6:CF:80:ARG:CG	6:CF:88:VAL:HB	2.44	0.47
7:CG:99:LEU:O	7:CG:100:ALA:C	2.52	0.47
10:CJ:33:GLN:O	10:CJ:75:ILE:HG23	2.13	0.47
11:CK:87:THR:HG22	11:CK:88:GLY:N	2.29	0.47
12:CL:54:LYS:N	12:CL:54:LYS:CD	2.75	0.47
13:CM:107:ALA:H	13:CM:108:ARG:HD2	1.79	0.47
13:CM:7:VAL:HG12	13:CM:7:VAL:O	2.14	0.47
19:CS:15:LEU:HD22	19:CS:15:LEU:N	2.28	0.47
19:CS:36:ARG:NH2	19:CS:72:GLY:HA2	2.29	0.47
23:CW:55:5MU:H73	23:CW:56:U:C4	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:CY:150:SER:C	25:CY:152:ASP:H	2.17	0.47
25:CY:3:LEU:CD1	25:CY:3:LEU:H	2.12	0.47
27:D1:13:ILE:HG13	27:D1:14:VAL:HG12	1.95	0.47
28:D2:17:SER:O	28:D2:20:GLU:N	2.47	0.47
28:D2:60:LEU:HG	28:D2:61:LEU:N	2.17	0.47
33:D7:29:LYS:HZ3	33:D7:32:LYS:HZ2	1.62	0.47
35:DA:942:G:H1'	35:DA:1189:A:H2	1.78	0.47
35:DA:1406:U:H3'	35:DA:1407:C:C6	2.47	0.47
35:DA:1459:G:H5''	35:DA:1460:A:OP2	2.14	0.47
35:DA:1707:G:C4	35:DA:1708:C:C5	3.02	0.47
35:DA:1767:C:O2	35:DA:1985:G:N2	2.44	0.47
35:DA:1917:U:H2'	35:DA:1918:A:H8	1.80	0.47
35:DA:2102:U:H2'	35:DA:2103:C:C5	2.49	0.47
35:DA:2354:G:O2'	35:DA:2355:C:H5'	2.14	0.47
35:DA:2850:A:C2'	35:DA:2851:A:H8	2.21	0.47
35:DA:778:G:H4'	38:DD:48:ARG:HD2	1.96	0.47
38:DD:118:VAL:HG22	38:DD:119:ALA:H	1.76	0.47
38:DD:117:VAL:HG21	38:DD:128:GLY:C	2.34	0.47
38:DD:270:ILE:O	38:DD:271:ILE:HG13	2.13	0.47
38:DD:46:GLN:N	38:DD:46:GLN:OE1	2.47	0.47
38:DD:94:LEU:CD1	38:DD:94:LEU:O	2.61	0.47
39:DE:179:GLU:O	39:DE:180:ASN:CB	2.63	0.47
39:DE:44:TYR:CD1	39:DE:44:TYR:N	2.83	0.47
39:DE:55:ASN:ND2	39:DE:75:VAL:HG22	2.28	0.47
40:DF:126:VAL:CG1	40:DF:193:VAL:HG13	2.45	0.47
40:DF:24:LEU:CD2	40:DF:24:LEU:N	2.78	0.47
41:DG:102:PHE:CD1	41:DG:106:LEU:HD22	2.49	0.47
41:DG:111:LEU:O	41:DG:114:ILE:HG12	2.14	0.47
42:DH:103:LEU:HD22	42:DH:123:PHE:CE2	2.49	0.47
44:DN:72:TYR:HB2	44:DN:85:ILE:HB	1.96	0.47
45:DO:63:VAL:HG22	45:DO:84:ALA:CA	2.32	0.47
45:DO:71:ARG:HG3	45:DO:71:ARG:NH1	2.28	0.47
35:DA:631:A:H4'	46:DP:65:ARG:HG3	1.96	0.47
46:DP:97:PRO:C	46:DP:99:LEU:N	2.68	0.47
48:DR:32:GLY:O	48:DR:116:LEU:HB2	2.14	0.47
48:DR:18:LEU:C	48:DR:18:LEU:HD13	2.34	0.47
35:DA:2838:G:N2	48:DR:93:GLY:HA3	2.29	0.47
50:DT:10:VAL:O	50:DT:11:GLU:C	2.52	0.47
50:DT:6:LEU:HD23	50:DT:6:LEU:O	2.14	0.47
55:DY:18:GLY:O	55:DY:20:TYR:N	2.46	0.47
55:DY:7:VAL:HG23	55:DY:8:LYS:HD2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1015:A:O5'	1:AA:1015:A:H8	1.98	0.47
1:AA:1426:C:C2	1:AA:1427:U:C5	3.02	0.47
1:AA:1527:C:C2	1:AA:1528:U:C5	3.02	0.47
1:AA:339:C:O2'	1:AA:340:U:H5'	2.14	0.47
1:AA:603:U:H2'	1:AA:604:G:C8	2.50	0.47
1:AA:769:G:O2'	1:AA:770:C:H5'	2.14	0.47
1:AA:811:C:H4'	1:AA:900:A:H62	1.79	0.47
1:AA:922:G:C6	1:AA:923:A:C6	3.02	0.47
1:AA:939:G:C4	1:AA:940:C:C5	3.02	0.47
2:AB:163:PHE:O	2:AB:164:VAL:HG23	2.14	0.47
2:AB:222:ILE:C	2:AB:222:ILE:HD13	2.34	0.47
2:AB:50:GLU:O	2:AB:51:LEU:C	2.53	0.47
2:AB:83:MET:O	2:AB:86:GLU:N	2.47	0.47
5:AE:147:ASP:HA	5:AE:150:ARG:CB	2.42	0.47
7:AG:102:ARG:O	7:AG:103:TRP:C	2.50	0.47
8:AH:60:ARG:HH11	8:AH:60:ARG:HG3	1.80	0.47
9:AI:4:TYR:HA	9:AI:88:TYR:CE1	2.49	0.47
11:AK:99:GLN:NE2	11:AK:105:VAL:HG11	2.29	0.47
14:AN:37:PHE:CE1	14:AN:53:LEU:HD22	2.49	0.47
20:AT:22:ARG:O	20:AT:23:ARG:C	2.53	0.47
21:AU:13:ILE:HA	21:AU:22:ARG:NH1	2.29	0.47
29:B3:17:LYS:HG2	35:BA:969:U:OP1	2.14	0.47
33:B7:34:ARG:NE	33:B7:39:ARG:HD2	2.28	0.47
34:B8:56:GLU:HA	34:B8:59:LYS:CE	2.44	0.47
35:BA:1166:C:H2'	35:BA:1167:U:C6	2.49	0.47
35:BA:1284:A:H2'	35:BA:1285:G:O4'	2.13	0.47
35:BA:1467:C:H2'	35:BA:1468:C:H6	1.79	0.47
35:BA:154(A):C:H5	35:BA:171:G:H1	1.59	0.47
35:BA:1555:G:H2'	35:BA:1556:C:H6	1.79	0.47
35:BA:1654:A:C2	39:BE:113:PHE:CD1	3.03	0.47
35:BA:2034:U:H2'	35:BA:2035:G:H5'	1.96	0.47
35:BA:2228:G:P	38:BD:263:ARG:HH21	2.37	0.47
35:BA:2519:U:OP1	35:BA:2519:U:H3'	2.14	0.47
35:BA:2701:C:H2'	35:BA:2702:U:H6	1.79	0.47
35:BA:2726:U:H6	45:BO:67:LYS:NZ	2.12	0.47
35:BA:604:G:H2'	35:BA:605:C:C6	2.49	0.47
35:BA:768:G:H2'	35:BA:769:G:C8	2.49	0.47
35:BA:828:U:O2	35:BA:828:U:H3'	2.14	0.47
35:BA:951:C:O2'	35:BA:952:G:H5'	2.14	0.47
35:BA:2127:G:H5'	37:BC:36:LYS:NZ	2.30	0.47
35:BA:2203:U:H4'	38:BD:151:LYS:HE3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:25:THR:O	38:BD:26:LYS:C	2.53	0.47
40:BF:103:LYS:O	40:BF:104:LYS:C	2.51	0.47
40:BF:119:ARG:HH11	40:BF:119:ARG:HG2	1.79	0.47
40:BF:202:PHE:CE1	40:BF:206:ILE:HD13	2.47	0.47
41:BG:161:THR:HG22	41:BG:163:ALA:H	1.78	0.47
42:BH:86:GLU:N	42:BH:86:GLU:OE1	2.47	0.47
44:BN:9:VAL:HG21	44:BN:39:ARG:HH21	1.79	0.47
46:BP:83:VAL:HG11	46:BP:112:LEU:HD21	1.96	0.47
49:BS:101:LEU:HD21	49:BS:103:GLU:CG	2.45	0.47
49:BS:92:TYR:HD1	49:BS:92:TYR:C	2.17	0.47
1:AA:1442(A):G:N2	50:BT:119:LYS:H	2.12	0.47
51:BU:98:LEU:C	51:BU:100:VAL:H	2.16	0.47
53:BW:24:ILE:O	53:BW:71:VAL:HG11	2.14	0.47
55:BY:37:VAL:CG2	55:BY:67:LEU:HG	2.43	0.47
55:BY:67:LEU:HD12	55:BY:68:HIS:N	2.29	0.47
56:BZ:66:SER:C	56:BZ:67:LEU:HD12	2.34	0.47
1:CA:1210:C:H4'	1:CA:1214:C:C4	2.49	0.47
1:CA:1271:G:H2'	1:CA:1272:G:O4'	2.14	0.47
1:CA:251:G:N2	1:CA:253:U:C4	2.83	0.47
1:CA:259:G:O2'	1:CA:260:G:H5'	2.13	0.47
1:CA:299:G:C6	1:CA:300:A:N1	2.83	0.47
1:CA:408:A:H2'	1:CA:409:G:O4'	2.15	0.47
5:CE:126:ARG:O	5:CE:128:PRO:N	2.47	0.47
5:CE:131:ILE:CD1	5:CE:131:ILE:H	2.21	0.47
8:CH:87:SER:HB3	8:CH:132:GLU:OE2	2.15	0.47
7:CG:16:LEU:CD1	9:CI:41:VAL:HG12	2.43	0.47
10:CJ:8:LEU:HB2	10:CJ:16:LEU:HD11	1.95	0.47
11:CK:88:GLY:N	11:CK:91:ARG:HB2	2.30	0.47
12:CL:41:ARG:HH11	12:CL:41:ARG:CB	2.19	0.47
14:CN:42:ILE:O	14:CN:45:ARG:N	2.48	0.47
15:CO:77:ARG:O	15:CO:80:ALA:HB3	2.14	0.47
17:CQ:56:VAL:HG23	17:CQ:78:GLU:HG3	1.96	0.47
17:CQ:62:SER:CB	17:CQ:72:ARG:HG3	2.44	0.47
20:CT:90:GLN:C	20:CT:93:GLU:OE2	2.52	0.47
28:D2:25:VAL:HG13	28:D2:26:ARG:N	2.29	0.47
32:D6:28:ARG:O	32:D6:29:ASN:C	2.51	0.47
34:D8:59:LYS:HB2	34:D8:59:LYS:NZ	2.29	0.47
35:DA:1037:G:H1	35:DA:1118:C:N4	2.06	0.47
35:DA:109:G:H2'	35:DA:110:G:H8	1.78	0.47
35:DA:1387:C:C5'	35:DA:1469:A:H4'	2.44	0.47
35:DA:1930:G:N2	35:DA:1968:G:C2'	2.74	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2115:G:H22	35:DA:2170:A:N6	2.12	0.47
26:D0:16:SER:HB3	35:DA:2262:U:OP2	2.14	0.47
35:DA:2404:C:N4	35:DA:2414:G:O6	2.47	0.47
35:DA:2495:G:C6	35:DA:2496:C:C4	3.02	0.47
35:DA:2556:C:H2'	35:DA:2557:G:O4'	2.13	0.47
35:DA:2763:G:H5'	35:DA:2763:G:C8	2.49	0.47
35:DA:2822:G:OP2	39:DE:110:GLY:O	2.32	0.47
35:DA:296:C:N4	35:DA:343:C:H42	2.12	0.47
35:DA:571:A:O2'	35:DA:573:G:O5'	2.32	0.47
35:DA:576:U:H2'	35:DA:577:G:H8	1.80	0.47
35:DA:619:G:O5'	35:DA:620:G:N2	2.43	0.47
34:D8:17:THR:HG22	35:DA:650:C:O2'	2.14	0.47
36:DB:73:A:H2'	36:DB:74:U:O4'	2.15	0.47
38:DD:167:GLY:O	38:DD:168:ARG:HG2	2.14	0.47
39:DE:143:ASN:N	39:DE:143:ASN:ND2	2.62	0.47
39:DE:4:ILE:HG12	39:DE:28:ALA:HB1	1.94	0.47
39:DE:60:ASN:O	39:DE:62:PRO:N	2.47	0.47
40:DF:28:ILE:HD13	40:DF:28:ILE:H	1.77	0.47
40:DF:43:LYS:HA	40:DF:98:SER:HA	1.95	0.47
41:DG:134:GLY:HA3	41:DG:155:MET:O	2.15	0.47
41:DG:55:LYS:O	41:DG:55:LYS:HD3	2.14	0.47
43:DI:20:ASP:O	43:DI:21:VAL:HG13	2.13	0.47
44:DN:40:PRO:CA	51:DU:64:ARG:HH22	2.27	0.47
45:DO:112:MET:O	45:DO:115:VAL:HG23	2.14	0.47
51:DU:108:GLU:HB3	51:DU:112:ARG:NH1	2.29	0.47
51:DU:29:SER:O	51:DU:30:LYS:NZ	2.43	0.47
51:DU:49:HIS:O	51:DU:52:ARG:N	2.48	0.47
51:DU:96:ALA:C	51:DU:98:LEU:H	2.18	0.47
52:DV:93:GLU:HG2	52:DV:94:LEU:N	2.28	0.47
53:DW:26:GLY:H	53:DW:71:VAL:CG1	2.27	0.47
55:DY:28:LYS:HZ2	55:DY:37:VAL:HA	1.77	0.47
56:DZ:117:LEU:HD23	56:DZ:117:LEU:H	1.79	0.47
56:DZ:150:LEU:O	56:DZ:171:ILE:HG12	2.13	0.47
1:AA:1149:C:H2'	1:AA:1150:U:C6	2.46	0.47
1:AA:1268:A:H2'	1:AA:1269:A:C8	2.48	0.47
1:AA:1273:G:H2'	1:AA:1274:G:O4'	2.14	0.47
1:AA:1442(B):A:N6	50:BT:118:ARG:NH2	2.63	0.47
1:AA:1443:G:N1	1:AA:1460:A:N3	2.62	0.47
1:AA:245:C:C2'	1:AA:246:A:H5'	2.44	0.47
1:AA:802:A:C2'	1:AA:803:G:H5'	2.43	0.47
2:AB:114:ARG:HA	2:AB:117:GLU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:213:LEU:C	2:AB:213:LEU:CD2	2.83	0.47
2:AB:69:LEU:C	2:AB:69:LEU:CD1	2.83	0.47
3:AC:172:ARG:NH1	3:AC:174:PRO:HG2	2.30	0.47
3:AC:40:ARG:O	3:AC:44:GLU:N	2.44	0.47
4:AD:100:ARG:O	4:AD:101:LEU:C	2.52	0.47
4:AD:206:PHE:O	4:AD:206:PHE:CG	2.68	0.47
4:AD:4:TYR:O	4:AD:5:ILE:CB	2.62	0.47
4:AD:5:ILE:CG2	4:AD:6:GLY:H	2.11	0.47
5:AE:127:ASN:O	5:AE:128:PRO:C	2.53	0.47
6:AF:10:LEU:HD12	6:AF:10:LEU:N	2.28	0.47
6:AF:72:VAL:CG1	6:AF:73:ASN:H	2.22	0.47
7:AG:32:ARG:O	7:AG:33:ASP:HB2	2.14	0.47
9:AI:7:THR:H	9:AI:83:ARG:HD2	1.80	0.47
12:AL:84:LEU:HD22	12:AL:101:VAL:HG21	1.97	0.47
13:AM:49:THR:HG22	13:AM:50:GLU:N	2.30	0.47
13:AM:50:GLU:O	13:AM:54:VAL:HG23	2.15	0.47
14:AN:42:ILE:O	14:AN:45:ARG:N	2.45	0.47
17:AQ:56:VAL:HG23	17:AQ:78:GLU:HG3	1.96	0.47
17:AQ:70:ARG:C	17:AQ:71:PHE:HD2	2.18	0.47
19:AS:15:LEU:N	19:AS:15:LEU:HD22	2.26	0.47
19:AS:72:GLY:C	19:AS:74:PHE:N	2.67	0.47
25:AY:76:LEU:O	25:AY:79:ILE:HB	2.13	0.47
25:AY:74:ASN:HA	25:AY:77:LYS:CG	2.44	0.47
27:B1:20:ARG:CZ	27:B1:41:ARG:NE	2.77	0.47
27:B1:85:LEU:CA	27:B1:87:PRO:HD3	2.43	0.47
27:B1:88:LYS:HG3	27:B1:89:GLU:OE2	2.15	0.47
32:B6:51:GLU:N	32:B6:51:GLU:OE1	2.45	0.47
34:B8:3:LYS:O	34:B8:4:MET:O	2.32	0.47
35:BA:1007:C:HO2'	44:BN:108:PRO:HA	1.77	0.47
35:BA:142:A:C8	35:BA:1408:C:H1'	2.49	0.47
35:BA:1659:U:OP2	39:BE:132:HIS:HE1	1.98	0.47
35:BA:1866:C:H2'	35:BA:1876:A:O4'	2.15	0.47
35:BA:1997:G:C2	35:BA:1998:G:C5	3.03	0.47
35:BA:2031:A:C6	35:BA:2498:C:H1'	2.50	0.47
35:BA:1136:G:N3	35:BA:2038:G:H4'	2.29	0.47
35:BA:2319:G:H4'	35:BA:2319:G:OP2	2.14	0.47
35:BA:2357:U:H2'	35:BA:2358:G:H5''	1.96	0.47
35:BA:2373:G:C6	35:BA:2374:C:N4	2.83	0.47
35:BA:2730:C:H2'	35:BA:2731:G:H8	1.80	0.47
35:BA:2843:G:C2	35:BA:2875:C:N3	2.82	0.47
35:BA:462:C:N4	35:BA:468:G:C6	2.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:70:G:H21	35:BA:71:A:N6	2.12	0.47
35:BA:957:A:H8	35:BA:957:A:OP1	1.98	0.47
36:BB:110:G:H2'	36:BB:111:G:H8	1.79	0.47
38:BD:5:LYS:N	38:BD:5:LYS:HD2	2.28	0.47
39:BE:181:LEU:CD2	39:BE:181:LEU:N	2.76	0.47
40:BF:153:SER:HA	40:BF:172:TRP:O	2.14	0.47
40:BF:185:ASP:HA	40:BF:188:ARG:CB	2.39	0.47
35:BA:660:G:C5'	40:BF:99:TYR:CD2	2.97	0.47
41:BG:107:LEU:HD11	41:BG:178:PHE:CE1	2.48	0.47
41:BG:174:GLU:O	41:BG:176:LEU:N	2.47	0.47
41:BG:72:ARG:HB3	41:BG:86:MET:N	2.19	0.47
35:BA:2667:C:H1'	42:BH:109:PHE:CE2	2.50	0.47
43:BI:29:TYR:CE1	43:BI:33:ARG:NE	2.82	0.47
43:BI:79:ILE:CG2	43:BI:81:VAL:HG23	2.45	0.47
45:BO:65:THR:O	45:BO:79:PHE:HB2	2.14	0.47
46:BP:84:ASN:CG	46:BP:116:GLY:HA3	2.34	0.47
35:BA:196:A:C5'	46:BP:46:LYS:HZ1	2.24	0.47
48:BR:56:LYS:HD2	48:BR:88:ARG:HA	1.97	0.47
49:BS:36:TYR:HA	49:BS:52:SER:HB2	1.96	0.47
49:BS:54:LEU:CD1	49:BS:58:LEU:O	2.61	0.47
49:BS:65:VAL:HG12	49:BS:69:VAL:HB	1.96	0.47
50:BT:27:THR:HG23	50:BT:28:VAL:N	2.28	0.47
50:BT:61:PHE:CZ	50:BT:76:PHE:HB3	2.49	0.47
50:BT:50:ILE:HD11	50:BT:64:ARG:HB3	1.96	0.47
52:BV:30:GLY:HA2	52:BV:64:HIS:O	2.14	0.47
52:BV:51:VAL:HG12	52:BV:52:VAL:N	2.30	0.47
54:BX:33:LYS:O	54:BX:35:THR:N	2.47	0.47
55:BY:2:ARG:HG2	55:BY:2:ARG:NH1	2.30	0.47
56:BZ:115:GLY:CA	56:BZ:175:VAL:O	2.51	0.47
1:CA:1039:C:H2'	1:CA:1040:U:C6	2.50	0.47
1:CA:1226:C:OP1	13:CM:96:LEU:HD13	2.15	0.47
1:CA:13:U:C5	1:CA:916:G:O6	2.68	0.47
1:CA:363:A:O2'	1:CA:364:A:H5'	2.13	0.47
1:CA:38:G:C2	1:CA:397:A:C2	3.02	0.47
1:CA:517:G:N3	1:CA:531:U:H5'	2.29	0.47
1:CA:630:G:H2'	1:CA:631:G:C5'	2.44	0.47
1:CA:788:U:H2'	1:CA:789:U:H6	1.79	0.47
2:CB:195:ASP:O	8:CH:68:ARG:NH2	2.47	0.47
3:CC:127:ARG:HD2	3:CC:127:ARG:N	2.29	0.47
4:CD:148:VAL:O	4:CD:149:ALA:O	2.33	0.47
4:CD:98:GLU:C	4:CD:100:ARG:H	2.17	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:105:VAL:HB	5:CE:106:PRO:HD3	1.97	0.47
6:CF:43:LEU:HD12	6:CF:43:LEU:N	2.26	0.47
6:CF:45:LEU:CD1	6:CF:46:ARG:N	2.76	0.47
7:CG:148:ASN:C	7:CG:150:ALA:N	2.68	0.47
7:CG:85:TYR:CE1	7:CG:154:TYR:HE1	2.33	0.47
9:CI:13:ALA:HB2	9:CI:68:GLY:HA3	1.96	0.47
1:CA:972:C:H4'	10:CJ:57:LYS:HG2	1.93	0.47
11:CK:73:MET:HA	11:CK:77:MET:HB3	1.97	0.47
1:CA:528:C:N4	12:CL:49:ASN:HD22	2.07	0.47
12:CL:55:VAL:CG1	12:CL:56:ALA:H	2.08	0.47
12:CL:91:LYS:HG3	12:CL:91:LYS:O	2.14	0.47
15:CO:87:ILE:HG22	15:CO:88:ARG:H	1.79	0.47
18:CR:81:PHE:O	18:CR:82:THR:CB	2.63	0.47
19:CS:16:LEU:O	19:CS:20:LEU:N	2.44	0.47
21:CU:12:LYS:HG2	21:CU:22:ARG:HB2	1.97	0.47
25:CY:177:GLU:O	25:CY:178:LYS:C	2.53	0.47
29:D3:32:GLN:OE1	29:D3:32:GLN:HA	2.14	0.47
34:D8:21:LYS:O	34:D8:23:VAL:HG23	2.14	0.47
34:D8:48:PHE:CD1	34:D8:48:PHE:N	2.76	0.47
35:DA:1022:G:O2'	35:DA:1023:U:OP2	2.26	0.47
35:DA:1112:G:O2'	35:DA:1113:U:H5''	2.15	0.47
35:DA:996:A:N6	35:DA:1160:G:C6	2.82	0.47
35:DA:814:C:H1'	35:DA:1225:G:H21	1.79	0.47
35:DA:1468:C:O2'	35:DA:1469:A:H5'	2.14	0.47
35:DA:1950:G:H8	35:DA:1950:G:O5'	1.98	0.47
35:DA:2001:A:C2	35:DA:2002:G:C4	3.02	0.47
35:DA:2003:G:C6	35:DA:2004:G:C5	3.02	0.47
35:DA:2319:G:C5	35:DA:2320:A:N6	2.83	0.47
35:DA:474:G:H4'	35:DA:475:U:OP1	2.12	0.47
35:DA:705:A:C2	35:DA:706:A:C4	3.02	0.47
35:DA:834:C:H2'	35:DA:835:A:C8	2.49	0.47
35:DA:979:G:N2	35:DA:985:C:N4	2.63	0.47
36:DB:40:U:H2'	36:DB:41:U:OP1	2.14	0.47
36:DB:61:G:H2'	36:DB:62:C:H6	1.78	0.47
36:DB:79:C:H2'	36:DB:80:U:O4'	2.15	0.47
38:DD:211:ARG:O	38:DD:212:SER:C	2.52	0.47
35:DA:1902:C:C4'	38:DD:244:ARG:HB2	2.43	0.47
40:DF:157:VAL:HB	40:DF:194:MET:HB3	1.97	0.47
41:DG:51:ARG:HE	41:DG:51:ARG:CA	2.21	0.47
42:DH:109:PHE:HE1	42:DH:152:ARG:NE	2.12	0.47
44:DN:123:TYR:N	44:DN:123:TYR:CD1	2.83	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:125:GLY:HA3	44:DN:126:PRO:O	2.14	0.47
44:DN:13:TRP:HD1	44:DN:13:TRP:H	1.58	0.47
45:DO:105:GLU:HA	45:DO:108:GLU:CD	2.34	0.47
46:DP:128:HIS:O	46:DP:129:ALA:HB2	2.14	0.47
47:DQ:35:VAL:HG22	47:DQ:100:GLY:O	2.14	0.47
48:DR:10:LEU:HD22	48:DR:17:ARG:CD	2.27	0.47
48:DR:17:ARG:CG	48:DR:17:ARG:HH11	2.23	0.47
50:DT:108:ARG:HB3	50:DT:108:ARG:HH11	1.80	0.47
52:DV:30:GLY:HA2	52:DV:64:HIS:O	2.14	0.47
52:DV:29:PRO:C	52:DV:31:ALA:H	2.18	0.47
55:DY:44:ILE:HG23	55:DY:45:VAL:N	2.30	0.47
56:DZ:125:LEU:CB	56:DZ:165:VAL:HG22	2.31	0.47
56:DZ:59:LEU:HD11	56:DZ:69:THR:OG1	2.15	0.47
56:DZ:94:GLU:HA	56:DZ:95:PRO:HD2	1.82	0.47
1:AA:1036:G:H2'	1:AA:1036:G:N3	2.28	0.47
1:AA:1300:G:O2'	1:AA:1301:U:O5'	2.26	0.47
1:AA:1321:C:H5''	1:AA:1322:C:H5''	1.92	0.47
1:AA:1463:C:H2'	1:AA:1464:G:C8	2.49	0.47
1:AA:1463:C:H2'	1:AA:1464:G:H8	1.80	0.47
1:AA:352:C:H2'	1:AA:352:C:O2	2.15	0.47
1:AA:729:A:H2'	1:AA:730:G:C8	2.49	0.47
1:AA:781:A:C2'	1:AA:782:A:H5'	2.44	0.47
1:AA:769:G:H1	1:AA:810:C:H42	1.61	0.47
2:AB:15:VAL:HG23	2:AB:209:ARG:HE	1.80	0.47
2:AB:224:GLN:HG3	2:AB:229:VAL:CG2	2.44	0.47
3:AC:105:GLU:CG	3:AC:106:VAL:H	2.14	0.47
4:AD:30:LYS:O	4:AD:32:ALA:N	2.47	0.47
5:AE:109:ILE:HD12	5:AE:135:THR:HB	1.97	0.47
5:AE:146:ALA:O	5:AE:148:VAL:N	2.48	0.47
7:AG:140:ASP:HA	7:AG:143:ARG:CZ	2.45	0.47
13:AM:107:ALA:H	13:AM:108:ARG:HD2	1.79	0.47
13:AM:36:LYS:O	13:AM:37:THR:HG23	2.15	0.47
13:AM:3:ARG:HG2	13:AM:9:ILE:HD11	1.95	0.47
13:AM:7:VAL:HG12	13:AM:7:VAL:O	2.15	0.47
15:AO:39:LEU:O	15:AO:42:HIS:HB3	2.14	0.47
19:AS:12:ASP:HB3	19:AS:15:LEU:CD2	2.44	0.47
23:AW:3:C:H6	23:AW:3:C:OP2	1.98	0.47
25:AY:116:ARG:HG2	25:AY:116:ARG:NH1	2.25	0.47
25:AY:150:SER:O	25:AY:152:ASP:N	2.47	0.47
26:B0:36:ILE:HA	26:B0:60:PHE:HB3	1.96	0.47
26:B0:77:ARG:HH22	35:BA:857:C:C5'	2.21	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:13:ILE:HB	27:B1:63:ALA:CB	2.41	0.47
29:B3:9:VAL:HG22	29:B3:54:VAL:HA	1.96	0.47
31:B5:40:LYS:HZ3	31:B5:45:VAL:CA	2.26	0.47
31:B5:51:TYR:C	31:B5:56:LYS:HG2	2.35	0.47
34:B8:2:PRO:C	34:B8:4:MET:H	2.17	0.47
34:B8:23:VAL:CG1	34:B8:46:ARG:NH1	2.76	0.47
34:B8:56:GLU:O	34:B8:57:ARG:C	2.52	0.47
35:BA:996:A:N6	35:BA:1160:G:C6	2.83	0.47
35:BA:1257:C:H2'	35:BA:1258:C:H6	1.80	0.47
35:BA:1317:A:C6	35:BA:1318:C:C4	3.03	0.47
35:BA:1916:A:H3'	35:BA:1917:U:C6	2.50	0.47
35:BA:199:A:C6	35:BA:2434:A:C6	3.02	0.47
35:BA:2596:U:H2'	35:BA:2597:G:O4'	2.15	0.47
35:BA:306:U:H2'	35:BA:306:U:O2	2.14	0.47
35:BA:38:A:H2'	35:BA:39:C:C6	2.49	0.47
35:BA:563:G:OP2	35:BA:572:A:H5'	2.15	0.47
36:BB:7:G:H2'	36:BB:8:U:O4'	2.15	0.47
35:BA:1902:C:C4'	38:BD:244:ARG:HB2	2.43	0.47
35:BA:2810:A:O2'	39:BE:61:ARG:NE	2.48	0.47
43:BI:102:SER:HB2	43:BI:109:ILE:CG2	2.44	0.47
44:BN:67:LEU:HB3	44:BN:88:GLU:CD	2.35	0.47
45:BO:111:PHE:CB	45:BO:114:ILE:HD13	2.35	0.47
45:BO:85:VAL:CG1	45:BO:86:ILE:N	2.77	0.47
46:BP:58:THR:C	46:BP:60:MET:N	2.66	0.47
35:BA:910:A:C4	47:BQ:13:GLN:OE1	2.68	0.47
49:BS:90:GLY:C	49:BS:92:TYR:N	2.63	0.47
51:BU:91:ASP:OD2	51:BU:96:ALA:HB2	2.15	0.47
56:BZ:54:HIS:HB3	56:BZ:101:PRO:CG	2.45	0.47
56:BZ:18:LEU:HD11	56:BZ:23:LYS:HB2	1.97	0.47
1:CA:1055:A:C2	1:CA:1056:U:H1'	2.50	0.47
1:CA:1238:A:N7	1:CA:1303:C:H1'	2.29	0.47
1:CA:1346:A:H5''	9:CI:120:ARG:NH1	2.24	0.47
1:CA:1392:G:O2'	1:CA:1393:U:H5'	2.13	0.47
1:CA:1457:G:H2'	1:CA:1458:G:C8	2.43	0.47
1:CA:1494:G:N3	1:CA:1494:G:H2'	2.30	0.47
1:CA:66:G:C4'	1:CA:173:U:C4	2.98	0.47
1:CA:66:G:C4'	1:CA:173:U:C5	2.93	0.47
1:CA:310:G:O2'	1:CA:311:C:H5'	2.15	0.47
1:CA:515:G:C2	1:CA:537:G:C2	3.03	0.47
1:CA:542:G:C4	1:CA:543:C:C5	3.03	0.47
1:CA:59:A:H5'	1:CA:60:A:H5'	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:774:G:H1	1:CA:805:C:N4	2.13	0.47
2:CB:187:LEU:HB2	2:CB:201:ILE:HB	1.96	0.47
2:CB:19:HIS:CE1	2:CB:206:ASP:HB2	2.50	0.47
2:CB:213:LEU:O	2:CB:217:ARG:HG2	2.14	0.47
2:CB:70:PHE:O	2:CB:92:TYR:HA	2.15	0.47
3:CC:33:LEU:O	3:CC:36:ASP:HB3	2.14	0.47
4:CD:13:ARG:NH1	4:CD:40:PRO:HA	2.29	0.47
4:CD:13:ARG:O	4:CD:15:GLU:N	2.47	0.47
4:CD:31:CYS:C	4:CD:33:MET:H	2.17	0.47
5:CE:6:PHE:HB2	5:CE:34:VAL:CG1	2.45	0.47
5:CE:78:HIS:HD2	8:CH:104:ARG:NE	2.12	0.47
8:CH:15:ASN:O	8:CH:16:ALA:C	2.53	0.47
8:CH:86:ILE:HG21	8:CH:133:LEU:CD2	2.44	0.47
9:CI:65:VAL:CG2	9:CI:66:ARG:H	2.27	0.47
10:CJ:34:VAL:HA	10:CJ:74:ILE:HA	1.97	0.47
11:CK:102:GLY:O	11:CK:103:LEU:O	2.33	0.47
11:CK:51:LYS:HA	11:CK:55:LYS:HG3	1.96	0.47
11:CK:58:PRO:HA	11:CK:90:GLY:HA3	1.96	0.47
13:CM:86:CYS:SG	13:CM:89:GLY:N	2.88	0.47
14:CN:6:LEU:O	14:CN:8:GLU:N	2.48	0.47
16:CP:4:ILE:HB	16:CP:66:PRO:CB	2.36	0.47
17:CQ:11:VAL:CA	17:CQ:53:LEU:HD11	2.44	0.47
19:CS:72:GLY:C	19:CS:74:PHE:N	2.68	0.47
26:D0:38:VAL:CG2	26:D0:59:LEU:HB2	2.45	0.47
27:D1:52:ARG:O	27:D1:53:VAL:HG12	2.15	0.47
27:D1:76:ARG:CA	27:D1:78:LYS:NZ	2.77	0.47
27:D1:87:PRO:HD2	27:D1:88:LYS:H	1.80	0.47
29:D3:52:HIS:CG	36:DB:83:G:H4'	2.48	0.47
29:D3:9:VAL:HG22	29:D3:54:VAL:HA	1.96	0.47
31:D5:30:LEU:HD23	31:D5:41:PRO:HB3	1.96	0.47
35:DA:1011:G:C6	35:DA:1151:G:C6	3.02	0.47
35:DA:1281:G:H1	35:DA:1286:A:N6	2.13	0.47
35:DA:1835:G:C5'	35:DA:1836:C:OP2	2.60	0.47
35:DA:1935:G:C3'	35:DA:1962:C:H42	2.27	0.47
31:D5:19:ARG:HA	35:DA:2046:G:O5'	2.14	0.47
35:DA:292:C:H2'	35:DA:292:C:O2	2.12	0.47
35:DA:419:C:O2'	35:DA:420:C:H5'	2.14	0.47
38:DD:243:GLY:O	38:DD:244:ARG:HB3	2.15	0.47
39:DE:107:THR:O	39:DE:190:GLY:HA2	2.15	0.47
40:DF:132:VAL:O	40:DF:133:ASN:C	2.53	0.47
35:DA:442:G:H4'	40:DF:46:ARG:HD3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:120:LEU:HD12	41:DG:180:PHE:CE1	2.50	0.47
41:DG:133:LEU:HB2	41:DG:134:GLY:H	1.54	0.47
44:DN:137:LYS:CG	44:DN:138:LEU:H	2.25	0.47
45:DO:14:THR:HG21	45:DO:86:ILE:HG12	1.96	0.47
45:DO:61:VAL:O	45:DO:84:ALA:CB	2.54	0.47
47:DQ:39:PRO:O	47:DQ:40:ALA:HB2	2.15	0.47
53:DW:20:VAL:O	53:DW:23:LEU:HB2	2.14	0.47
53:DW:28:SER:O	53:DW:29:LEU:C	2.52	0.47
54:DX:18:TYR:O	54:DX:21:PHE:HB2	2.14	0.47
54:DX:24:GLY:HA2	54:DX:80:ILE:HG13	1.95	0.47
54:DX:31:HIS:O	54:DX:32:PRO:C	2.53	0.47
1:AA:1351:U:O5'	1:AA:1351:U:H6	1.98	0.47
1:AA:237:C:C5'	17:AQ:25:ARG:NH1	2.77	0.47
1:AA:279:A:OP2	17:AQ:95:TYR:OH	2.30	0.47
1:AA:357:G:O2'	1:AA:358:U:H5'	2.15	0.47
1:AA:513:C:H2'	1:AA:514:C:C6	2.49	0.47
1:AA:689:C:H2'	1:AA:689:C:O2	2.13	0.47
2:AB:233:SER:C	2:AB:235:SER:H	2.18	0.47
3:AC:127:ARG:HD2	3:AC:127:ARG:N	2.30	0.47
3:AC:94:LEU:HD12	3:AC:95:THR:N	2.29	0.47
4:AD:147:ALA:HA	4:AD:182:LYS:HB3	1.96	0.47
4:AD:2:GLY:O	4:AD:3:ARG:C	2.52	0.47
4:AD:32:ALA:O	4:AD:36:ARG:N	2.48	0.47
4:AD:65:ARG:NH1	4:AD:72:GLU:HB2	2.30	0.47
5:AE:39:GLY:O	5:AE:69:VAL:N	2.48	0.47
6:AF:38:GLU:HG3	6:AF:65:VAL:HA	1.97	0.47
7:AG:119:ARG:O	7:AG:122:HIS:N	2.47	0.47
8:AH:31:PHE:HA	8:AH:34:GLU:HG2	1.97	0.47
9:AI:83:ARG:C	9:AI:86:VAL:HG12	2.35	0.47
1:AA:1125:U:O4	10:AJ:5:ARG:HD3	2.14	0.47
10:AJ:34:VAL:HA	10:AJ:74:ILE:HA	1.96	0.47
12:AL:89:ARG:HB2	12:AL:89:ARG:CZ	2.43	0.47
15:AO:83:GLU:O	15:AO:83:GLU:HG2	2.15	0.47
16:AP:2:VAL:O	16:AP:2:VAL:HG22	2.14	0.47
17:AQ:12:SER:HA	17:AQ:14:LYS:NZ	2.30	0.47
19:AS:17:GLU:C	19:AS:19:VAL:H	2.17	0.47
19:AS:36:ARG:NH2	19:AS:72:GLY:HA2	2.30	0.47
1:AA:1400:C:H5'	24:AX:18:C:N4	2.30	0.47
25:AY:107:THR:HG23	25:AY:110:ARG:H	1.79	0.47
25:AY:16:LYS:O	25:AY:19:GLU:N	2.48	0.47
25:AY:36:ALA:HA	25:AY:39:LEU:CD2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:14:VAL:O	27:B1:15:ALA:O	2.33	0.47
27:B1:71:TYR:O	27:B1:74:VAL:HB	2.15	0.47
34:B8:35:GLN:HE21	34:B8:36:LYS:HZ3	1.60	0.47
35:BA:1576:U:H2'	35:BA:1577:C:H6	1.79	0.47
1:AA:1494:G:H5''	35:BA:1913:A:N6	2.29	0.47
35:BA:2050:C:H2'	35:BA:2051:A:O4'	2.15	0.47
35:BA:2168:G:H2'	35:BA:2170:A:OP2	2.14	0.47
26:B0:17:GLN:HG2	35:BA:2261:C:P	2.54	0.47
35:BA:2274:A:N6	35:BA:2276:G:C4	2.83	0.47
35:BA:2426:A:C3'	35:BA:2427:C:C5'	2.86	0.47
35:BA:1938:A:N1	35:BA:2590:A:H1'	2.28	0.47
35:BA:301:G:H1'	35:BA:302:C:C6	2.49	0.47
35:BA:304:G:H2'	35:BA:305:U:C6	2.50	0.47
35:BA:669:G:N2	35:BA:670:A:C4	2.83	0.47
35:BA:751:A:H62	35:BA:789:A:N6	2.13	0.47
35:BA:798:G:C6	35:BA:799:G:C6	3.02	0.47
38:BD:12:SER:HB2	38:BD:208:LYS:HB3	1.96	0.47
35:BA:2051:A:C4'	39:BE:141:ILE:HD11	2.45	0.47
41:BG:108:ASN:O	41:BG:112:PRO:HG2	2.14	0.47
41:BG:104:GLU:O	41:BG:108:ASN:OD1	2.31	0.47
43:BI:92:VAL:O	43:BI:92:VAL:HG13	2.14	0.47
46:BP:35:HIS:CD2	46:BP:35:HIS:O	2.68	0.47
47:BQ:77:LYS:O	47:BQ:79:LEU:N	2.39	0.47
49:BS:89:ARG:HB3	49:BS:97:ARG:HH22	1.80	0.47
1:AA:1442(A):G:N2	50:BT:118:ARG:HB2	2.30	0.47
50:BT:61:PHE:CZ	50:BT:76:PHE:CB	2.97	0.47
51:BU:24:TYR:HB2	51:BU:29:SER:OG	2.14	0.47
54:BX:30:VAL:O	54:BX:31:HIS:C	2.52	0.47
54:BX:58:HIS:C	54:BX:59:VAL:HG22	2.35	0.47
55:BY:2:ARG:HH11	55:BY:2:ARG:HG2	1.78	0.47
56:BZ:102:LEU:HD12	56:BZ:121:HIS:O	2.14	0.47
1:CA:122:G:H2'	1:CA:123:C:C6	2.49	0.47
1:CA:189(D):C:H1'	1:CA:189(H):G:N2	2.30	0.47
1:CA:563:A:N3	1:CA:563:A:H2'	2.29	0.47
1:CA:77:G:H1	1:CA:92:C:N4	2.11	0.47
2:CB:12:GLU:C	2:CB:14:GLY:H	2.17	0.47
2:CB:222:ILE:HD13	2:CB:222:ILE:C	2.34	0.47
2:CB:212:GLN:NE2	2:CB:235:SER:HB3	2.30	0.47
3:CC:107:GLN:N	3:CC:107:GLN:CD	2.66	0.47
3:CC:207:VAL:O	3:CC:207:VAL:HG12	2.14	0.47
3:CC:40:ARG:O	3:CC:44:GLU:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:120:LEU:HB2	4:CD:126:ILE:HD11	1.96	0.47
4:CD:15:GLU:H	4:CD:15:GLU:CD	2.16	0.47
4:CD:187:ARG:NH1	4:CD:187:ARG:HG2	2.29	0.47
4:CD:80:GLU:HA	4:CD:80:GLU:OE2	2.15	0.47
7:CG:81:GLY:C	7:CG:83:ALA:H	2.18	0.47
8:CH:122:ARG:CA	8:CH:125:ARG:HB3	2.40	0.47
11:CK:30:VAL:CG2	11:CK:68:ALA:HB2	2.43	0.47
12:CL:41:ARG:NH1	12:CL:41:ARG:HB3	2.18	0.47
12:CL:66:VAL:HG21	12:CL:98:TYR:CD1	2.49	0.47
17:CQ:56:VAL:O	17:CQ:76:LEU:HD12	2.15	0.47
20:CT:83:ARG:CA	20:CT:86:ARG:HB3	2.39	0.47
22:CV:41:C:H2'	22:CV:42:C:H6	1.74	0.47
25:CY:25:LEU:C	25:CY:179:LYS:HE2	2.35	0.47
26:D0:37:LEU:C	26:D0:38:VAL:CG2	2.83	0.47
34:D8:39:LYS:HZ1	34:D8:43:GLN:HG3	1.80	0.47
35:DA:1110:G:H4'	35:DA:1110:G:OP1	2.15	0.47
35:DA:1180:C:C2'	35:DA:1181:C:H5'	2.45	0.47
35:DA:1279:G:H2'	35:DA:1280:G:H8	1.80	0.47
35:DA:1578:U:H6	35:DA:1578:U:OP2	1.97	0.47
35:DA:1578:U:H2'	35:DA:1579:A:H5''	1.96	0.47
35:DA:1622:G:C2	35:DA:1623:G:C8	3.02	0.47
35:DA:1910:G:O2'	35:DA:1911:U:H5'	2.15	0.47
35:DA:2357:U:H2'	35:DA:2358:G:H5''	1.96	0.47
35:DA:2531:A:O2'	35:DA:2532:G:H5'	2.13	0.47
35:DA:2571:C:H5'	35:DA:2572:A:C5'	2.45	0.47
35:DA:2606:C:O2'	35:DA:2607:G:H5'	2.14	0.47
35:DA:2712:U:C1'	35:DA:2712(A):A:C8	2.97	0.47
35:DA:675:A:C6	35:DA:676:A:C6	3.02	0.47
35:DA:680:G:O2'	35:DA:681:G:H5'	2.14	0.47
35:DA:79:G:H2'	35:DA:80:G:H8	1.79	0.47
35:DA:902:C:H2'	35:DA:903:C:H6	1.78	0.47
36:DB:28:C:H2'	36:DB:29:A:O4'	2.14	0.47
36:DB:29:A:OP2	49:DS:32:LEU:HB2	2.15	0.47
38:DD:267:SER:HA	38:DD:270:ILE:HG13	1.96	0.47
38:DD:25:THR:HG23	38:DD:27:THR:HB	1.96	0.47
40:DF:24:LEU:O	40:DF:25:PRO:C	2.52	0.47
41:DG:88:ILE:CG2	41:DG:89:GLY:N	2.77	0.47
42:DH:37:VAL:HG12	42:DH:38:SER:N	2.30	0.47
42:DH:86:GLU:N	42:DH:86:GLU:OE1	2.46	0.47
46:DP:23:PRO:HB2	46:DP:33:ARG:CG	2.43	0.47
47:DQ:16:ARG:HB3	47:DQ:16:ARG:CZ	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:101:LEU:HD21	49:DS:103:GLU:CG	2.45	0.47
49:DS:29:PHE:C	49:DS:29:PHE:CD1	2.87	0.47
50:DT:50:ILE:HA	50:DT:99:LEU:CD1	2.43	0.47
51:DU:31:SER:HB3	51:DU:34:LYS:CB	2.44	0.47
1:AA:1065:U:H4'	1:AA:1066:C:O5'	2.14	0.47
1:AA:1346:A:O3'	1:AA:1347:G:H4'	2.14	0.47
1:AA:269:C:H2'	1:AA:270:A:C8	2.50	0.47
1:AA:341:C:H2'	1:AA:342:C:H6	1.80	0.47
1:AA:515:G:C2	1:AA:537:G:C2	3.03	0.47
1:AA:739:C:O2'	1:AA:740:U:H5'	2.15	0.47
1:AA:77:G:H1	1:AA:92:C:N4	2.11	0.47
2:AB:187:LEU:HD23	2:AB:201:ILE:O	2.14	0.47
2:AB:221:LEU:O	2:AB:221:LEU:HD22	2.15	0.47
3:AC:29:TYR:HE2	14:AN:37:PHE:CE1	2.32	0.47
4:AD:176:LEU:HD12	4:AD:177:ASP:H	1.78	0.47
4:AD:13:ARG:HH22	4:AD:36:ARG:HH11	1.63	0.47
5:AE:136:MET:O	5:AE:137:GLU:C	2.53	0.47
6:AF:42:GLU:O	6:AF:44:GLY:N	2.48	0.47
9:AI:11:LYS:C	9:AI:13:ALA:N	2.68	0.47
9:AI:28:VAL:CG2	9:AI:63:ILE:HB	2.44	0.47
11:AK:19:ALA:CB	11:AK:32:ILE:HG12	2.45	0.47
11:AK:33:THR:HG22	11:AK:39:PRO:HA	1.96	0.47
13:AM:83:ASP:OD1	13:AM:86:CYS:HB2	2.14	0.47
15:AO:36:ILE:O	15:AO:37:ASN:C	2.53	0.47
15:AO:54:ARG:HG2	15:AO:54:ARG:NH1	2.29	0.47
18:AR:53:ARG:NH1	18:AR:58:LEU:O	2.48	0.47
18:AR:85:LEU:CG	18:AR:86:VAL:H	2.19	0.47
20:AT:16:HIS:O	20:AT:19:SER:N	2.47	0.47
25:AY:156:ARG:NH2	47:BQ:80:GLU:HG2	2.30	0.47
26:B0:20:ARG:CD	26:B0:20:ARG:N	2.78	0.47
28:B2:14:ARG:O	28:B2:15:LYS:C	2.53	0.47
28:B2:22:GLU:CA	28:B2:25:VAL:HG12	2.44	0.47
29:B3:13:ILE:HD12	35:BA:989:G:N7	2.30	0.47
29:B3:7:LYS:O	29:B3:9:VAL:HG13	2.14	0.47
31:B5:25:LEU:HD12	53:BW:23:LEU:HD11	1.94	0.47
35:BA:105:C:C5'	35:BA:106:C:OP2	2.62	0.47
35:BA:1257:C:H2'	35:BA:1258:C:C6	2.50	0.47
35:BA:1676:A:H2'	35:BA:1677:A:O4'	2.14	0.47
35:BA:1775:U:H2'	35:BA:1776:G:O5'	2.14	0.47
35:BA:1917:U:H2'	35:BA:1918:A:H8	1.79	0.47
35:BA:1982:C:H2'	35:BA:1983:C:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2533:A:C3'	35:BA:2534:A:H5''	2.45	0.47
35:BA:31:C:H2'	35:BA:32:C:O4'	2.14	0.47
35:BA:42:G:H2'	35:BA:43:A:H8	1.78	0.47
35:BA:481:G:H2'	35:BA:507:A:N1	2.28	0.47
35:BA:571:A:O2'	35:BA:573:G:O5'	2.32	0.47
35:BA:613:G:C6	35:BA:615:G:O6	2.68	0.47
35:BA:679:C:O2	35:BA:679:C:H2'	2.14	0.47
35:BA:717:G:H2'	35:BA:718:A:O4'	2.15	0.47
35:BA:814:C:HO2'	35:BA:815:C:H5'	1.79	0.47
35:BA:911:A:C4	47:BQ:9:TYR:OH	2.64	0.47
36:BB:28:C:H2'	36:BB:29:A:C8	2.50	0.47
38:BD:27:THR:O	38:BD:28:GLU:CB	2.61	0.47
39:BE:120:TRP:CD1	39:BE:155:LYS:HB3	2.49	0.47
35:BA:443:A:OP1	40:BF:46:ARG:CB	2.63	0.47
40:BF:71:GLY:O	40:BF:73:ALA:N	2.48	0.47
41:BG:111:LEU:O	41:BG:114:ILE:HG13	2.15	0.47
42:BH:78:GLY:O	42:BH:136:ILE:HG23	2.15	0.47
35:BA:195:A:OP1	46:BP:46:LYS:HE2	2.15	0.47
49:BS:66:ALA:HA	49:BS:69:VAL:CG1	2.44	0.47
51:BU:61:TRP:O	51:BU:63:VAL:N	2.48	0.47
51:BU:96:ALA:C	51:BU:98:LEU:N	2.68	0.47
52:BV:72:VAL:O	52:BV:73:SER:CB	2.62	0.47
35:BA:2012:G:O3'	53:BW:96:ILE:HD11	2.14	0.47
54:BX:12:VAL:HG13	54:BX:17:ALA:HB1	1.94	0.47
54:BX:36:LYS:O	54:BX:37:THR:C	2.52	0.47
54:BX:57:LEU:O	54:BX:58:HIS:CG	2.67	0.47
56:BZ:149:SER:CB	56:BZ:173:ALA:CA	2.90	0.47
1:CA:1296:C:H4'	1:CA:1302:U:O4	2.15	0.47
1:CA:1316:G:H2'	1:CA:1317:C:C5'	2.45	0.47
1:CA:1375:A:C4	1:CA:1376:U:C5	3.02	0.47
1:CA:598:U:H4'	8:CH:94:TYR:CD1	2.49	0.47
1:CA:929:G:C6	1:CA:930:C:C4	3.02	0.47
2:CB:177:ALA:O	2:CB:180:LEU:N	2.48	0.47
2:CB:36:ARG:HB3	2:CB:41:ILE:CD1	2.42	0.47
4:CD:100:ARG:O	4:CD:103:ASN:HB3	2.14	0.47
4:CD:102:ASP:HA	4:CD:121:VAL:HG21	1.96	0.47
7:CG:27:ILE:HD12	7:CG:27:ILE:H	1.80	0.47
13:CM:114:ARG:C	13:CM:116:THR:H	2.17	0.47
13:CM:27:LYS:O	13:CM:30:ALA:HB3	2.15	0.47
1:CA:658:G:O4'	15:CO:22:THR:HB	2.14	0.47
15:CO:53:HIS:O	15:CO:56:LEU:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CO:80:ALA:O	15:CO:81:LEU:C	2.53	0.47
23:CW:27:G:H2'	23:CW:28:U:O4'	2.14	0.47
25:CY:127:VAL:O	25:CY:130:ARG:N	2.48	0.47
25:CY:164:ILE:O	25:CY:168:PHE:HB2	2.15	0.47
25:CY:2:THR:HG23	25:CY:4:LYS:HB3	1.96	0.47
26:D0:36:ILE:O	26:D0:36:ILE:HG13	2.15	0.47
27:D1:87:PRO:CB	27:D1:91:LYS:NZ	2.72	0.47
34:D8:49:VAL:HG12	34:D8:53:PRO:HD3	1.97	0.47
35:DA:118:A:OP2	35:DA:119:A:H5''	2.15	0.47
35:DA:1578:U:C6	35:DA:1578:U:OP2	2.68	0.47
35:DA:1778:U:H2'	35:DA:1784:A:N6	2.30	0.47
35:DA:2016:U:H2'	35:DA:2017:U:C6	2.50	0.47
35:DA:2033:A:HO2'	35:DA:2034:U:P	2.36	0.47
35:DA:2079:U:H2'	35:DA:2080:G:H8	1.78	0.47
35:DA:2206:G:N2	35:DA:2207:G:C5'	2.72	0.47
35:DA:2536:G:C6	35:DA:2537:U:C4	3.02	0.47
35:DA:31:C:H2'	35:DA:32:C:O4'	2.14	0.47
35:DA:366:C:H5''	35:DA:403:U:H3	1.79	0.47
35:DA:393:C:H2'	35:DA:394:A:C8	2.45	0.47
35:DA:958:U:OP2	47:DQ:14:ARG:NH1	2.48	0.47
29:D3:17:LYS:HG2	35:DA:969:U:OP1	2.15	0.47
36:DB:28:C:H2'	36:DB:29:A:C8	2.49	0.47
38:DD:113:VAL:C	38:DD:115:GLN:H	2.14	0.47
38:DD:25:THR:CG2	38:DD:82:ILE:N	2.77	0.47
38:DD:27:THR:O	38:DD:28:GLU:CB	2.58	0.47
40:DF:119:ARG:HH11	40:DF:119:ARG:HG2	1.79	0.47
40:DF:197:ASP:O	40:DF:200:GLU:HB3	2.14	0.47
40:DF:45:ARG:HD2	40:DF:46:ARG:H	1.79	0.47
40:DF:89:VAL:O	40:DF:91:GLY:N	2.42	0.47
42:DH:89:ILE:O	42:DH:161:GLY:O	2.32	0.47
42:DH:25:LYS:HB3	42:DH:32:GLU:OE2	2.15	0.47
43:DI:123:LEU:HD11	43:DI:143:SER:O	2.15	0.47
43:DI:6:LEU:O	43:DI:8:PRO:N	2.47	0.47
46:DP:58:THR:C	46:DP:60:MET:N	2.66	0.47
49:DS:66:ALA:HB1	49:DS:98:VAL:O	2.15	0.47
49:DS:25:ARG:NH2	49:DS:89:ARG:NH1	2.58	0.47
50:DT:61:PHE:CZ	50:DT:76:PHE:CB	2.98	0.47
51:DU:20:LEU:H	51:DU:20:LEU:CD2	2.20	0.47
51:DU:35:ALA:O	51:DU:36:ARG:C	2.52	0.47
52:DV:34:GLU:HG2	52:DV:35:LEU:H	1.78	0.47
53:DW:17:VAL:O	53:DW:20:VAL:CG2	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:14:LEU:HD12	55:DY:23:ARG:O	2.15	0.47
55:DY:86:ARG:HG2	55:DY:87:LYS:H	1.80	0.47
56:DZ:44:PHE:O	56:DZ:48:PHE:N	2.44	0.47
1:AA:1108:G:H2'	1:AA:1108:G:N3	2.30	0.47
1:AA:1109:C:O2'	1:AA:1110:A:H5'	2.14	0.47
1:AA:951:G:C6	1:AA:1231:G:C6	3.02	0.47
1:AA:1238:A:N7	1:AA:1303:C:H1'	2.30	0.47
1:AA:1306:A:H2'	1:AA:1307:U:O4'	2.15	0.47
1:AA:334:C:C2'	1:AA:335:C:H5'	2.45	0.47
1:AA:39:G:H2'	1:AA:40:C:H6	1.80	0.47
1:AA:509:A:H4'	1:AA:510:A:OP1	2.15	0.47
1:AA:564:C:C5	17:AQ:31:LEU:HD11	2.49	0.47
1:AA:59:A:C5'	1:AA:60:A:C5'	2.85	0.47
1:AA:629:G:H2'	1:AA:630:G:O4'	2.14	0.47
2:AB:85:ALA:HB1	2:AB:92:TYR:HB3	1.96	0.47
4:AD:31:CYS:C	4:AD:33:MET:H	2.18	0.47
4:AD:97:LEU:O	4:AD:97:LEU:CD2	2.63	0.47
7:AG:156:TRP:H	7:AG:156:TRP:HD1	1.63	0.47
8:AH:87:SER:HB3	8:AH:132:GLU:OE2	2.14	0.47
9:AI:48:GLU:N	9:AI:49:PRO:CD	2.76	0.47
9:AI:93:ARG:C	9:AI:95:LYS:H	2.18	0.47
9:AI:97:LYS:HB3	9:AI:98:PRO:HD3	1.96	0.47
1:AA:1060:C:H4'	10:AJ:52:GLY:N	2.30	0.47
11:AK:34:ASP:HB2	11:AK:35:PRO:CD	2.44	0.47
12:AL:22:SER:C	12:AL:24:VAL:N	2.68	0.47
23:AW:75:C:C5	23:AW:76:C:N3	2.82	0.47
25:AY:63:PRO:CB	25:AY:64:ARG:NH2	2.77	0.47
28:B2:50:ILE:HG23	28:B2:54:LYS:HZ3	1.80	0.47
35:BA:1279:G:H2'	35:BA:1280:G:C8	2.49	0.47
35:BA:1289:C:O2'	35:BA:1290:C:H5'	2.15	0.47
35:BA:1363:C:H2'	35:BA:1364:G:H8	1.80	0.47
35:BA:2200:C:N4	35:BA:2223:G:H1	2.08	0.47
35:BA:2292:C:H2'	35:BA:2293:C:C6	2.50	0.47
35:BA:2513:G:H2'	35:BA:2514:U:C6	2.49	0.47
35:BA:2663:G:O2'	35:BA:2664:G:H5'	2.14	0.47
35:BA:694:U:H2'	35:BA:695:G:O5'	2.14	0.47
35:BA:751:A:H62	35:BA:789:A:H62	1.63	0.47
35:BA:79:G:H2'	35:BA:80:G:H8	1.80	0.47
35:BA:809:G:H2'	35:BA:810:U:C6	2.50	0.47
35:BA:836:G:C5	35:BA:837:C:N4	2.83	0.47
38:BD:247:ALA:CB	38:BD:253:GLN:HA	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:119:ARG:HD2	39:BE:120:TRP:CE2	2.50	0.47
39:BE:8:LYS:HE3	39:BE:188:VAL:HG13	1.95	0.47
41:BG:100:TRP:HA	41:BG:100:TRP:CE3	2.50	0.47
42:BH:84:SER:O	42:BH:85:LYS:CB	2.60	0.47
43:BI:109:ILE:HD13	43:BI:111:PRO:HD3	1.96	0.47
43:BI:83:ALA:O	43:BI:144:VAL:HG13	2.15	0.47
35:BA:2094:G:P	43:BI:22:LYS:HD2	2.55	0.47
35:BA:1138:G:H1'	44:BN:105:GLY:O	2.14	0.47
35:BA:1952:A:P	45:BO:44:LYS:HZ3	2.37	0.47
45:BO:87:ILE:HG13	45:BO:91:LEU:CD1	2.45	0.47
35:BA:814:C:C5	46:BP:27:HIS:CE1	3.03	0.47
48:BR:55:ALA:CB	48:BR:79:LEU:HD12	2.43	0.47
49:BS:26:LEU:HD13	49:BS:87:PHE:HD1	1.80	0.47
53:BW:86:LEU:HD12	53:BW:87:PRO:CD	2.45	0.47
1:CA:1015:A:H2'	1:CA:1016:A:H8	1.79	0.47
1:CA:122:G:H2'	1:CA:123:C:H6	1.79	0.47
1:CA:1383:C:O2'	1:CA:1384:C:H5'	2.14	0.47
1:CA:1446:U:H4'	1:CA:1447:A:C6	2.50	0.47
1:CA:1499:A:O2'	1:CA:1500:A:H5'	2.14	0.47
1:CA:27:G:H2'	1:CA:28:G:H8	1.79	0.47
1:CA:44:G:N2	1:CA:45:U:H1'	2.30	0.47
1:CA:596:C:O2	1:CA:596:C:H2'	2.14	0.47
1:CA:596:C:O2	1:CA:597:G:C8	2.67	0.47
1:CA:929:G:C2'	1:CA:930:C:H5'	2.44	0.47
2:CB:193:ASP:O	2:CB:194:PRO:O	2.32	0.47
2:CB:41:ILE:N	2:CB:41:ILE:CD1	2.78	0.47
4:CD:114:ARG:O	4:CD:117:ALA:HB3	2.15	0.47
5:CE:39:GLY:O	5:CE:69:VAL:N	2.48	0.47
7:CG:115:ARG:HB2	7:CG:118:VAL:HG21	1.96	0.47
9:CI:126:SER:O	9:CI:127:LYS:CB	2.60	0.47
13:CM:74:VAL:HA	13:CM:77:ASN:HD22	1.80	0.47
15:CO:11:VAL:HG21	15:CO:34:LEU:CD2	2.45	0.47
17:CQ:34:LYS:O	17:CQ:36:ILE:HG12	2.14	0.47
18:CR:36:ASN:HD21	18:CR:40:LEU:HD21	1.80	0.47
20:CT:100:ILE:CD1	20:CT:100:ILE:N	2.70	0.47
35:DA:1175:U:C4'	35:DA:1176:G:H3'	2.45	0.47
35:DA:1448:G:H5'	35:DA:1449:A:OP1	2.14	0.47
35:DA:1496:A:H2'	35:DA:1498:C:H5	1.79	0.47
35:DA:1686:C:H3'	35:DA:1687:G:H8	1.80	0.47
35:DA:1778:U:C5	35:DA:1784:A:C2	3.03	0.47
35:DA:2070:G:C2	35:DA:2071:A:C4	3.02	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2242:G:H2'	35:DA:2243:U:O4'	2.15	0.47
35:DA:2747:G:C2	35:DA:2754:U:C4	3.03	0.47
35:DA:2752:C:C5	35:DA:2753:A:N7	2.82	0.47
35:DA:562:U:O2'	35:DA:563:G:H5''	2.14	0.47
35:DA:845:G:OP2	35:DA:845:G:H8	1.97	0.47
38:DD:62:TYR:OH	38:DD:64:ILE:HD12	2.15	0.47
39:DE:19:ARG:O	39:DE:19:ARG:HG3	2.14	0.47
39:DE:13:ARG:HA	39:DE:22:PRO:HA	1.97	0.47
40:DF:12:LEU:O	40:DF:14:PRO:HD3	2.15	0.47
42:DH:13:LYS:CA	42:DH:13:LYS:HE2	2.36	0.47
42:DH:30:LYS:NZ	42:DH:81:GLU:HA	2.30	0.47
44:DN:67:LEU:C	44:DN:69:GLN:N	2.68	0.47
45:DO:23:ARG:HG3	45:DO:24:VAL:N	2.30	0.47
47:DQ:35:VAL:CG2	47:DQ:101:ARG:O	2.62	0.47
48:DR:87:TYR:OH	48:DR:116:LEU:HD22	2.14	0.47
49:DS:101:LEU:O	49:DS:101:LEU:HD13	2.13	0.47
35:DA:2334:G:C5'	49:DS:13:ARG:HG2	2.34	0.47
50:DT:110:ILE:CG2	50:DT:111:ARG:N	2.78	0.47
51:DU:74:LEU:HD11	51:DU:79:PHE:HB2	1.96	0.47
54:DX:35:THR:HG23	54:DX:36:LYS:N	2.30	0.47
55:DY:38:ILE:CG2	55:DY:39:VAL:N	2.77	0.47
56:DZ:126:VAL:HA	56:DZ:164:ALA:HB3	1.97	0.47
56:DZ:29:TYR:O	56:DZ:30:ASN:HB3	2.15	0.47
1:AA:1226:C:H5'	19:AS:80:TYR:CE2	2.50	0.47
1:AA:1445:C:H6	1:AA:1445:C:O5'	1.98	0.47
1:AA:149:A:C2	1:AA:150:C:C4	3.03	0.47
1:AA:408:A:C2	1:AA:409:G:C4	3.03	0.47
1:AA:407:G:N1	1:AA:408:A:C5	2.83	0.47
1:AA:596:C:O2	1:AA:597:G:C8	2.68	0.47
1:AA:741:G:H5'	15:AO:39:LEU:HD21	1.96	0.47
3:AC:155:GLY:HA3	3:AC:163:ALA:CB	2.43	0.47
3:AC:178:LEU:C	3:AC:180:ALA:H	2.17	0.47
4:AD:4:TYR:O	4:AD:5:ILE:HB	2.14	0.47
4:AD:60:GLU:O	4:AD:61:LYS:C	2.54	0.47
6:AF:27:GLN:O	6:AF:31:GLU:HG3	2.15	0.47
7:AG:122:HIS:O	7:AG:125:MET:HB3	2.14	0.47
8:AH:35:ILE:N	8:AH:35:ILE:HD13	2.30	0.47
8:AH:95:VAL:HG23	8:AH:95:VAL:O	2.15	0.47
9:AI:37:PHE:CZ	9:AI:74:ILE:HG12	2.49	0.47
13:AM:10:PRO:CB	13:AM:18:ALA:HB1	2.44	0.47
15:AO:39:LEU:HD12	15:AO:56:LEU:HD12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:60:ILE:HG12	17:AQ:61:GLU:N	2.30	0.47
17:AQ:62:SER:OG	17:AQ:72:ARG:HG3	2.15	0.47
19:AS:6:LYS:CD	19:AS:6:LYS:N	2.78	0.47
26:B0:11:ARG:O	26:B0:11:ARG:HG2	2.15	0.47
27:B1:87:PRO:CD	27:B1:88:LYS:H	2.27	0.47
28:B2:12:GLU:O	28:B2:12:GLU:OE2	2.32	0.47
34:B8:4:MET:HG3	34:B8:4:MET:O	2.15	0.47
35:BA:1210:A:H1'	35:BA:1212:G:C4	2.50	0.47
35:BA:1271:G:C2	35:BA:1617:C:H4'	2.50	0.47
35:BA:1441:G:O2'	35:BA:1442:G:H5'	2.15	0.47
35:BA:2019:A:N1	35:BA:2020:A:C4	2.82	0.47
35:BA:2171:A:C8	35:BA:2172:U:C5	3.03	0.47
35:BA:2366:A:H2'	35:BA:2367:G:O4'	2.15	0.47
35:BA:2758:A:C3'	35:BA:2759:G:H5''	2.44	0.47
35:BA:626:U:H5''	35:BA:627:A:H5'	1.96	0.47
35:BA:802:A:H2'	35:BA:802:A:N3	2.29	0.47
37:BC:76:ALA:C	37:BC:78:ALA:H	2.16	0.47
38:BD:213:ARG:O	38:BD:215:LEU:N	2.48	0.47
38:BD:85:ASP:OD1	38:BD:87:ASN:HB2	2.15	0.47
39:BE:132:HIS:CD2	39:BE:135:HIS:NE2	2.82	0.47
40:BF:170:LEU:HD23	40:BF:173:VAL:HB	1.97	0.47
40:BF:45:ARG:CD	40:BF:46:ARG:H	2.28	0.47
40:BF:89:VAL:CG1	40:BF:90:PHE:H	2.23	0.47
41:BG:85:GLY:C	41:BG:87:PRO:HD3	2.35	0.47
42:BH:89:ILE:O	42:BH:161:GLY:O	2.33	0.47
42:BH:17:VAL:HG12	42:BH:19:VAL:HG23	1.97	0.47
44:BN:17:ASP:OD2	44:BN:17:ASP:O	2.33	0.47
45:BO:105:GLU:HA	45:BO:108:GLU:CD	2.35	0.47
34:B8:30:ARG:NE	46:BP:62:LEU:HB2	2.30	0.47
47:BQ:73:PRO:HG3	47:BQ:93:TYR:CE2	2.50	0.47
49:BS:38:GLN:CD	49:BS:47:THR:HG23	2.35	0.47
52:BV:78:LYS:HD3	52:BV:79:VAL:CA	2.44	0.47
52:BV:90:PRO:CD	52:BV:91:TYR:H	2.26	0.47
53:BW:10:VAL:HG23	53:BW:101:SER:O	2.15	0.47
55:BY:18:GLY:C	55:BY:20:TYR:N	2.66	0.47
55:BY:86:ARG:HG2	55:BY:87:LYS:H	1.78	0.47
1:CA:355:C:O2'	1:CA:356:A:H5'	2.15	0.47
1:CA:552:U:H2'	1:CA:553:A:C8	2.50	0.47
1:CA:603:U:H2'	1:CA:604:G:C8	2.50	0.47
1:CA:611:A:O2'	1:CA:612:C:H5'	2.15	0.47
1:CA:625:G:O2'	1:CA:626:U:H5'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:961:U:O2'	1:CA:962:C:H5'	2.15	0.47
1:CA:972:C:OP2	10:CJ:57:LYS:HE2	2.15	0.47
4:CD:101:LEU:HD23	4:CD:121:VAL:CG1	2.44	0.47
4:CD:42:GLN:OE1	4:CD:42:GLN:C	2.54	0.47
8:CH:47:GLY:O	8:CH:62:TYR:N	2.39	0.47
9:CI:27:THR:O	9:CI:28:VAL:CG2	2.63	0.47
13:CM:90:LEU:O	13:CM:91:ARG:HB2	2.15	0.47
15:CO:47:LYS:H	15:CO:47:LYS:HD3	1.80	0.47
19:CS:63:THR:HG22	19:CS:66:MET:CE	2.45	0.47
20:CT:83:ARG:HA	20:CT:86:ARG:HD3	1.96	0.47
21:CU:13:ILE:HA	21:CU:22:ARG:NH1	2.30	0.47
25:CY:120:GLN:O	25:CY:121:TYR:C	2.53	0.47
25:CY:15:GLN:N	25:CY:168:PHE:CZ	2.82	0.47
27:D1:25:LYS:O	27:D1:26:ARG:O	2.33	0.47
27:D1:83:GLU:CG	27:D1:86:SER:CB	2.92	0.47
29:D3:45:GLY:HA3	35:DA:852:G:H5'	1.95	0.47
35:DA:1006:C:H2'	35:DA:1007:C:H6	1.79	0.47
35:DA:1048:A:N6	35:DA:1106:A:C8	2.83	0.47
35:DA:1424:G:OP1	38:DD:33:LEU:HD21	2.15	0.47
35:DA:1860:G:H2'	35:DA:1861:G:C8	2.50	0.47
35:DA:2205:C:C2	35:DA:2220:G:N1	2.83	0.47
35:DA:2383:G:H2'	35:DA:2384:G:C8	2.50	0.47
35:DA:2487:G:O2'	35:DA:2488:A:H5'	2.15	0.47
35:DA:2497:A:O5'	35:DA:2497:A:C8	2.68	0.47
35:DA:252:G:OP2	46:DP:50:ARG:NH1	2.47	0.47
35:DA:2517:C:C6	35:DA:2542:A:C2	3.03	0.47
35:DA:2831:G:O4'	35:DA:2883:A:C2	2.68	0.47
35:DA:315:G:H2'	35:DA:316:C:C6	2.50	0.47
35:DA:314:A:O2'	35:DA:315:G:H5'	2.14	0.47
35:DA:447:A:C4	35:DA:473:G:N7	2.83	0.47
35:DA:744:G:OP1	39:DE:132:HIS:CB	2.63	0.47
37:DC:124:GLY:O	37:DC:125:SER:CB	2.63	0.47
38:DD:211:ARG:HH11	38:DD:211:ARG:HG2	1.79	0.47
38:DD:31:LYS:O	38:DD:32:SER:O	2.33	0.47
39:DE:47:VAL:HG12	39:DE:49:LEU:HD11	1.97	0.47
40:DF:28:ILE:N	40:DF:28:ILE:CD1	2.78	0.47
41:DG:60:LEU:HA	41:DG:63:ILE:HG12	1.96	0.47
42:DH:157:TYR:CD1	42:DH:170:ARG:O	2.68	0.47
43:DI:88:ILE:HG13	43:DI:122:GLU:HA	1.96	0.47
45:DO:81:ASP:CG	45:DO:81:ASP:O	2.53	0.47
47:DQ:32:TYR:HA	47:DQ:132:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:88:GLY:O	47:DQ:89:ASN:CB	2.61	0.47
48:DR:29:LEU:HG	48:DR:79:LEU:HD23	1.97	0.47
49:DS:19:LYS:HB3	49:DS:20:ARG:NH2	2.30	0.47
49:DS:87:PHE:HZ	49:DS:97:ARG:HH21	1.61	0.47
50:DT:26:ASP:O	50:DT:88:ILE:HB	2.15	0.47
50:DT:31:SER:O	50:DT:32:TYR:O	2.33	0.47
50:DT:33:LYS:HE2	50:DT:43:GLN:OE1	2.15	0.47
50:DT:61:PHE:CE2	50:DT:76:PHE:HB3	2.50	0.47
51:DU:111:GLU:O	51:DU:113:ALA:N	2.48	0.47
35:DA:1341:U:H5'	54:DX:57:LEU:CD2	2.45	0.47
55:DY:2:ARG:HH11	55:DY:2:ARG:HG2	1.80	0.47
55:DY:28:LYS:CE	55:DY:37:VAL:HA	2.45	0.47
35:DA:300:A:H5''	55:DY:97:ARG:NH1	2.30	0.47
56:DZ:100:VAL:O	56:DZ:101:PRO:O	2.31	0.47
56:DZ:10:ARG:HB2	56:DZ:36:LYS:O	2.15	0.47
1:AA:1419:G:N2	1:AA:1482:G:C1'	2.78	0.47
1:AA:1428:A:H2'	1:AA:1429:C:H6	1.73	0.47
1:AA:375:U:OP1	16:AP:69:THR:HG21	2.14	0.47
1:AA:376:G:O3'	16:AP:5:ARG:HD2	2.15	0.47
1:AA:577:G:C8	1:AA:816:A:C6	3.02	0.47
1:AA:93:G:C6	1:AA:96:U:C4	3.03	0.47
2:AB:99:GLY:O	2:AB:100:GLY:C	2.54	0.47
3:AC:182:ILE:HG12	3:AC:203:PHE:HD1	1.79	0.47
6:AF:37:VAL:HG13	6:AF:65:VAL:HG11	1.97	0.47
1:AA:823:G:H21	8:AH:1:MET:HE1	1.79	0.47
9:AI:71:SER:O	9:AI:74:ILE:HB	2.14	0.47
15:AO:70:LEU:CD2	15:AO:78:TYR:HA	2.44	0.47
15:AO:87:ILE:HG22	15:AO:88:ARG:H	1.80	0.47
27:B1:66:HIS:O	27:B1:69:LYS:HB2	2.15	0.47
28:B2:21:LEU:O	28:B2:25:VAL:HG12	2.14	0.47
28:B2:57:ILE:HD11	28:B2:59:ARG:CZ	2.45	0.47
29:B3:23:LEU:N	29:B3:23:LEU:HD12	2.30	0.47
29:B3:25:ALA:HB3	29:B3:26:LEU:HD23	1.97	0.47
33:B7:33:ARG:O	33:B7:36:GLN:HB3	2.14	0.47
35:BA:1163:G:C2	35:BA:1164:G:C8	3.03	0.47
35:BA:1364:G:N3	35:BA:1364:G:H2'	2.30	0.47
35:BA:1371:G:H8	35:BA:1371:G:O5'	1.98	0.47
35:BA:1668:A:N3	35:BA:1670:C:C4	2.83	0.47
35:BA:2165:G:H3'	35:BA:2166:G:H8	1.79	0.47
35:BA:2415:G:C2	35:BA:2416:C:C2	3.03	0.47
35:BA:271(W):G:H5'	35:BA:271(X):G:OP2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2870:C:H2'	35:BA:2871:C:C5'	2.45	0.47
35:BA:8:A:H2	35:BA:2896:C:O2	1.97	0.47
35:BA:49:A:N6	35:BA:177:G:C5	2.82	0.47
35:BA:701:G:C2'	35:BA:702:G:H5'	2.45	0.47
35:BA:997:G:OP1	51:BU:93:LYS:HE3	2.15	0.47
35:BA:2572:A:OP2	39:BE:144:ARG:HB2	2.14	0.47
40:BF:155:LEU:HD13	40:BF:174:VAL:HB	1.96	0.47
40:BF:45:ARG:CG	40:BF:46:ARG:N	2.78	0.47
41:BG:86:MET:O	41:BG:87:PRO:O	2.32	0.47
43:BI:119:PRO:O	43:BI:121:LYS:N	2.48	0.47
44:BN:128:HIS:HD2	44:BN:131:GLN:CB	2.28	0.47
44:BN:84:LYS:O	44:BN:85:ILE:HD13	2.15	0.47
45:BO:14:THR:HG21	45:BO:86:ILE:HG12	1.96	0.47
35:BA:2482:G:H1	47:BQ:53:ALA:HB2	1.80	0.47
48:BR:94:TYR:CD1	48:BR:94:TYR:N	2.82	0.47
50:BT:45:PHE:CZ	50:BT:74:ARG:HB2	2.49	0.47
50:BT:88:ILE:HG22	50:BT:89:VAL:HG23	1.97	0.47
53:BW:5:ALA:HB3	53:BW:105:VAL:N	2.29	0.47
54:BX:14:SER:O	54:BX:15:GLU:C	2.53	0.47
54:BX:26:TYR:N	54:BX:26:TYR:CD1	2.83	0.47
54:BX:62:LYS:HD2	54:BX:68:ARG:HD2	1.97	0.47
54:BX:75:ASP:C	54:BX:76:ARG:HG3	2.35	0.47
55:BY:86:ARG:CB	55:BY:88:LYS:HZ2	2.22	0.47
56:BZ:9:TYR:CE2	56:BZ:35:ARG:HB3	2.50	0.47
1:CA:1223:C:P	19:CS:78:ARG:HH12	2.38	0.47
1:CA:1513:A:C2	1:CA:1514:C:C4	3.03	0.47
1:CA:1524:C:C2	1:CA:1525:G:C8	3.03	0.47
1:CA:66:G:N2	1:CA:172:A:C2	2.83	0.47
1:CA:221:C:O2'	1:CA:222:U:H5'	2.15	0.47
1:CA:241:C:H1'	1:CA:286:G:N2	2.29	0.47
1:CA:522:C:H1'	1:CA:536:C:H5'	1.97	0.47
1:CA:661:G:H1	1:CA:744:C:H42	1.62	0.47
1:CA:779:C:O2'	11:CK:120:ARG:HD3	2.15	0.47
2:CB:157:ARG:HG2	2:CB:158:LEU:N	2.27	0.47
5:CE:82:VAL:CG2	5:CE:138:ALA:HA	2.39	0.47
6:CF:40:VAL:HG23	6:CF:62:TRP:O	2.15	0.47
6:CF:71:ARG:O	6:CF:72:VAL:C	2.53	0.47
9:CI:50:LEU:HB3	9:CI:55:ALA:CB	2.45	0.47
10:CJ:94:VAL:CG1	10:CJ:95:GLU:N	2.77	0.47
11:CK:19:ALA:CB	11:CK:32:ILE:HG12	2.45	0.47
7:CG:149:ARG:HD3	11:CK:59:TYR:CZ	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CM:28:ALA:C	13:CM:30:ALA:N	2.69	0.47
21:CU:5:ASP:OD1	21:CU:6:ARG:N	2.47	0.47
25:CY:160:GLU:O	25:CY:163:LYS:HB3	2.15	0.47
25:CY:165:THR:O	25:CY:166:ASP:O	2.33	0.47
25:CY:2:THR:HG23	25:CY:2:THR:O	2.15	0.47
26:D0:12:ASN:O	26:D0:14:ARG:N	2.47	0.47
27:D1:83:GLU:HG2	27:D1:86:SER:CB	2.43	0.47
35:DA:1011:G:C5	35:DA:1151:G:N1	2.82	0.47
35:DA:1027:A:O2'	35:DA:1028:A:H5'	2.15	0.47
35:DA:1223:G:C5	35:DA:1225:G:OP2	2.68	0.47
35:DA:1257:C:C2	35:DA:1258:C:C5	3.03	0.47
35:DA:1317:A:C6	35:DA:1318:C:C4	3.03	0.47
35:DA:151:C:H2'	35:DA:152:G:C8	2.49	0.47
35:DA:1300:U:H1'	35:DA:1626:G:N2	2.30	0.47
35:DA:1853:A:H1'	35:DA:2233:U:O2'	2.15	0.47
35:DA:1975:G:C6	35:DA:1976:U:C4	3.03	0.47
35:DA:1998:G:H4'	35:DA:2724:C:H4'	1.97	0.47
35:DA:2031:A:C6	35:DA:2498:C:H1'	2.50	0.47
35:DA:2041:U:H2'	35:DA:2042:A:C8	2.49	0.47
35:DA:2168:G:H2'	35:DA:2170:A:OP2	2.15	0.47
35:DA:412:A:N6	35:DA:2411:A:H2'	2.29	0.47
35:DA:2472:G:H5'	35:DA:2473:U:C5'	2.45	0.47
35:DA:926:A:H2'	35:DA:927:G:C8	2.45	0.47
35:DA:996:A:H2'	35:DA:997:G:C8	2.47	0.47
36:DB:64:C:H2'	36:DB:65:C:H6	1.79	0.47
38:DD:106:ILE:HD13	38:DD:157:ARG:HB2	1.97	0.47
39:DE:103:ASP:HA	39:DE:168:MET:HA	1.96	0.47
39:DE:47:VAL:CG1	39:DE:49:LEU:HD11	2.44	0.47
40:DF:71:GLY:O	40:DF:73:ALA:N	2.48	0.47
41:DG:63:ILE:HB	41:DG:141:PHE:CD1	2.50	0.47
41:DG:57:ALA:O	41:DG:61:ALA:HB2	2.15	0.47
41:DG:91:ARG:HD2	41:DG:92:VAL:CA	2.44	0.47
42:DH:18:GLU:HB3	42:DH:25:LYS:HD2	1.96	0.47
43:DI:29:TYR:C	43:DI:32:PRO:HD2	2.35	0.47
43:DI:2:LYS:O	43:DI:39:ALA:HB2	2.15	0.47
43:DI:58:LEU:HD23	43:DI:58:LEU:O	2.15	0.47
44:DN:115:ARG:O	44:DN:118:LYS:N	2.47	0.47
44:DN:56:ASN:CA	44:DN:125:GLY:H	2.26	0.47
44:DN:28:THR:HG22	44:DN:29:LYS:N	2.30	0.47
44:DN:29:LYS:C	44:DN:31:ALA:N	2.68	0.47
45:DO:113:LYS:O	45:DO:114:ILE:C	2.51	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:114:ILE:HD11	46:DP:130:PHE:CE2	2.50	0.47
46:DP:135:LEU:CD1	46:DP:139:LYS:HD2	2.44	0.47
46:DP:77:ARG:HB2	46:DP:78:PRO:CD	2.37	0.47
47:DQ:74:TYR:HB3	47:DQ:91:GLU:OE2	2.14	0.47
48:DR:101:ALA:O	48:DR:102:GLU:HB2	2.15	0.47
48:DR:97:VAL:HG13	48:DR:114:VAL:HG23	1.96	0.47
49:DS:15:ARG:HA	49:DS:17:ARG:HG2	1.97	0.47
49:DS:17:ARG:NE	49:DS:89:ARG:NH2	2.63	0.47
49:DS:66:ALA:CA	49:DS:69:VAL:HG12	2.45	0.47
49:DS:90:GLY:C	49:DS:92:TYR:H	2.15	0.47
53:DW:14:PRO:HG3	53:DW:101:SER:OG	2.15	0.47
55:DY:77:PRO:O	55:DY:78:ALA:CB	2.64	0.47
56:DZ:163:LEU:HD23	56:DZ:163:LEU:N	2.20	0.47
56:DZ:51:ALA:O	56:DZ:52:SER:OG	2.29	0.47
1:AA:1184:G:C2	1:AA:1185:G:C5	3.03	0.46
1:AA:678:U:H3	1:AA:713:G:N2	2.12	0.46
1:AA:885:G:H1	1:AA:912:C:N4	2.13	0.46
3:AC:130:VAL:HG21	3:AC:157:ILE:HG23	1.97	0.46
3:AC:41:GLY:O	3:AC:45:LYS:HG3	2.15	0.46
4:AD:98:GLU:C	4:AD:100:ARG:H	2.18	0.46
4:AD:64:LEU:HD23	4:AD:203:VAL:HG21	1.97	0.46
4:AD:80:GLU:HA	4:AD:80:GLU:OE2	2.15	0.46
6:AF:87:ARG:NH1	6:AF:87:ARG:HG3	2.30	0.46
8:AH:83:ILE:CD1	8:AH:137:VAL:HG22	2.38	0.46
12:AL:58:VAL:CG1	12:AL:59:ARG:N	2.77	0.46
14:AN:22:THR:HB	14:AN:33:VAL:HG21	1.97	0.46
14:AN:39:LEU:CD1	14:AN:47:LEU:HD12	2.44	0.46
15:AO:41:GLU:O	15:AO:44:LYS:HB3	2.15	0.46
27:B1:83:GLU:HG3	27:B1:85:LEU:HB2	1.97	0.46
29:B3:48:GLU:O	29:B3:51:ALA:HB2	2.15	0.46
35:BA:565:C:H4'	35:BA:1253:A:N6	2.30	0.46
35:BA:1317:A:C2	35:BA:1318:C:C2	3.03	0.46
35:BA:1339:G:N2	35:BA:1340:U:H5	2.12	0.46
35:BA:1450:G:H1	35:BA:1462:C:H42	1.63	0.46
35:BA:1494:A:C3'	35:BA:1494:A:N3	2.76	0.46
35:BA:1609:A:C5	35:BA:1616:A:C8	3.03	0.46
35:BA:1638:C:H2'	35:BA:1639:U:O4'	2.16	0.46
35:BA:1793:C:O2'	35:BA:1794:U:H5'	2.15	0.46
35:BA:1991:U:H2'	35:BA:1992:G:H5''	1.98	0.46
35:BA:528:A:C2	35:BA:2043:C:H5'	2.33	0.46
35:BA:2088:G:H2'	35:BA:2089:U:O4'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2230:G:O2'	35:BA:2231:C:H5'	2.15	0.46
35:BA:2419:U:O2'	35:BA:2420:C:H5'	2.14	0.46
35:BA:2531:A:O2'	35:BA:2532:G:H5'	2.15	0.46
35:BA:2778:A:H4'	35:BA:2779:U:OP2	2.14	0.46
35:BA:2824:C:H2'	35:BA:2825:C:O4'	2.14	0.46
35:BA:287:C:H2'	35:BA:288:C:O4'	2.16	0.46
35:BA:297:C:H2'	35:BA:298:G:O4'	2.14	0.46
35:BA:348:G:H2'	35:BA:349:G:H8	1.80	0.46
35:BA:464:U:H2'	35:BA:465:G:C8	2.50	0.46
35:BA:776:G:O6	35:BA:793:A:H2'	2.16	0.46
35:BA:810:U:C2	46:BP:31:ALA:O	2.68	0.46
39:BE:120:TRP:O	39:BE:122:PHE:N	2.47	0.46
41:BG:130:ASN:HA	41:BG:161:THR:HB	1.96	0.46
41:BG:42:GLY:HA2	41:BG:89:GLY:CA	2.46	0.46
41:BG:58:GLN:HG3	41:BG:59:GLU:N	2.30	0.46
42:BH:19:VAL:HG22	42:BH:24:VAL:HG13	1.96	0.46
44:BN:122:VAL:CG1	44:BN:123:TYR:N	2.77	0.46
47:BQ:41:TRP:HB3	47:BQ:94:VAL:HG21	1.97	0.46
50:BT:11:GLU:OE1	50:BT:11:GLU:N	2.47	0.46
51:BU:31:SER:C	51:BU:33:ARG:N	2.68	0.46
52:BV:83:ARG:CG	52:BV:83:ARG:NH1	2.65	0.46
52:BV:93:GLU:HG2	52:BV:94:LEU:H	1.79	0.46
53:BW:65:LEU:CD2	53:BW:67:ASP:HB2	2.45	0.46
54:BX:31:HIS:CG	54:BX:32:PRO:HD2	2.50	0.46
35:BA:143:G:C4'	54:BX:38:GLU:HG3	2.45	0.46
1:CA:1054:C:O2	1:CA:1054:C:H2'	2.14	0.46
1:CA:1255:G:O2'	1:CA:1258:G:H1'	2.15	0.46
1:CA:1303:C:O2	1:CA:1303:C:C2'	2.60	0.46
1:CA:322:C:O2'	1:CA:323:U:H5'	2.15	0.46
1:CA:450:G:H4'	16:CP:41:PRO:O	2.15	0.46
1:CA:885:G:H2'	1:CA:886:G:C8	2.51	0.46
2:CB:114:ARG:HA	2:CB:117:GLU:HB2	1.96	0.46
4:CD:17:VAL:CG1	4:CD:18:LYS:H	2.26	0.46
7:CG:42:ILE:HG23	7:CG:117:ALA:CB	2.45	0.46
7:CG:115:ARG:O	7:CG:118:VAL:HG23	2.15	0.46
7:CG:50:ILE:O	7:CG:50:ILE:HG22	2.15	0.46
8:CH:110:ALA:HA	8:CH:136:GLU:HA	1.96	0.46
1:CA:1060:C:H4'	10:CJ:52:GLY:N	2.30	0.46
1:CA:502:G:OP1	12:CL:118:SER:N	2.48	0.46
13:CM:22:ILE:HD12	13:CM:22:ILE:N	2.30	0.46
16:CP:20:VAL:HG21	16:CP:32:TYR:CG	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:30:ASP:O	18:CR:32:ARG:N	2.45	0.46
18:CR:53:ARG:NH1	18:CR:58:LEU:O	2.48	0.46
18:CR:66:LEU:O	18:CR:70:ILE:HG13	2.16	0.46
25:CY:92:PRO:HA	25:CY:100:TYR:O	2.15	0.46
25:CY:107:THR:HA	25:CY:111:ARG:HH12	1.80	0.46
25:CY:46:TYR:C	25:CY:48:ALA:N	2.68	0.46
26:D0:40:GLN:O	26:D0:57:PHE:HB2	2.15	0.46
26:D0:29:GLN:HB2	26:D0:67:VAL:HG21	1.97	0.46
27:D1:62:VAL:CG2	27:D1:66:HIS:O	2.62	0.46
27:D1:88:LYS:HB2	27:D1:92:LYS:HD3	1.97	0.46
33:D7:31:LEU:HD23	33:D7:42:LEU:HD22	1.97	0.46
33:D7:34:ARG:HE	33:D7:39:ARG:NE	2.12	0.46
34:D8:39:LYS:CD	34:D8:39:LYS:C	2.84	0.46
35:DA:1493:C:O2	35:DA:1493:C:C2'	2.63	0.46
35:DA:1349:A:N6	35:DA:1598:C:N4	2.63	0.46
35:DA:1747:G:O2'	35:DA:1747(A):G:H5'	2.14	0.46
35:DA:185:U:O2'	35:DA:186:G:H5'	2.16	0.46
35:DA:1885:A:H2'	35:DA:1886:C:O4'	2.15	0.46
35:DA:210:C:H2'	35:DA:211:A:H8	1.80	0.46
35:DA:2597:G:H2'	35:DA:2598:A:C8	2.51	0.46
35:DA:2631:G:H22	39:DE:61:ARG:NH1	2.11	0.46
35:DA:2810:A:C2'	39:DE:61:ARG:NH2	2.78	0.46
35:DA:287:C:H2'	35:DA:288:C:O4'	2.15	0.46
35:DA:432:A:C6	35:DA:433:C:C4	3.02	0.46
35:DA:573:G:O2'	35:DA:574:C:H3'	2.15	0.46
35:DA:701:G:C2'	35:DA:702:G:H5'	2.45	0.46
35:DA:60:G:H21	35:DA:74:A:H2'	1.80	0.46
35:DA:791:C:O2	35:DA:794:G:H5'	2.15	0.46
37:DC:75:LEU:HA	37:DC:94:VAL:HG22	1.96	0.46
38:DD:223:GLY:C	38:DD:225:ALA:H	2.17	0.46
35:DA:1568:G:H4'	38:DD:59:LYS:HG3	1.97	0.46
39:DE:110:GLY:HA2	39:DE:161:GLY:CA	2.26	0.46
40:DF:121:GLY:O	40:DF:123:LEU:N	2.48	0.46
40:DF:25:PRO:HB3	40:DF:119:ARG:HD3	1.96	0.46
41:DG:141:PHE:C	41:DG:144:ILE:HG22	2.36	0.46
41:DG:141:PHE:O	41:DG:144:ILE:HG22	2.15	0.46
42:DH:144:VAL:CG1	42:DH:148:ILE:HD11	2.43	0.46
42:DH:38:SER:C	42:DH:40:GLU:H	2.18	0.46
42:DH:61:HIS:C	42:DH:63:SER:H	2.18	0.46
43:DI:115:ALA:HB2	43:DI:129:THR:O	2.15	0.46
43:DI:85:GLU:O	43:DI:86:THR:O	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:42:TRP:CH2	44:DN:44:PRO:HA	2.50	0.46
44:DN:95:PRO:O	44:DN:98:VAL:HG23	2.15	0.46
45:DO:25:LEU:HD13	45:DO:38:VAL:HG11	1.96	0.46
46:DP:75:ILE:HD13	46:DP:77:ARG:NH2	2.30	0.46
47:DQ:77:LYS:C	47:DQ:79:LEU:H	2.19	0.46
48:DR:26:LYS:NZ	48:DR:71:GLN:HB3	2.30	0.46
49:DS:36:TYR:HA	49:DS:52:SER:HB2	1.96	0.46
35:DA:993:G:H1'	52:DV:91:TYR:CE1	2.50	0.46
53:DW:29:LEU:HD21	53:DW:33:ARG:CZ	2.43	0.46
54:DX:49:VAL:HG13	54:DX:85:PRO:HB3	1.97	0.46
55:DY:8:LYS:N	55:DY:8:LYS:HD2	2.20	0.46
1:AA:1015:A:H2'	1:AA:1016:A:C8	2.50	0.46
1:AA:1057:G:C2'	1:AA:1058:G:H5'	2.45	0.46
1:AA:1246:C:H2'	1:AA:1247:U:H6	1.80	0.46
1:AA:1304:G:N7	1:AA:1305:G:C6	2.83	0.46
1:AA:1380:U:O2	7:AG:3:ARG:NH1	2.48	0.46
1:AA:1433:A:O2'	1:AA:1434:A:H5'	2.15	0.46
1:AA:1422:G:N2	1:AA:1479:C:C2	2.83	0.46
1:AA:221:C:O2'	1:AA:222:U:H5'	2.15	0.46
1:AA:241:C:O2'	1:AA:242:C:H5'	2.14	0.46
1:AA:332:G:O2'	1:AA:333:G:H5'	2.15	0.46
1:AA:32:A:C2	1:AA:33:A:C4	3.03	0.46
1:AA:402:G:O2'	1:AA:403:C:H5'	2.15	0.46
1:AA:435:C:O2	1:AA:435:C:H2'	2.14	0.46
1:AA:473:G:H5''	16:AP:81:ARG:NE	2.30	0.46
1:AA:55:A:N7	1:AA:56:U:C5	2.84	0.46
2:AB:54:THR:HG21	2:AB:201:ILE:CD1	2.44	0.46
3:AC:186:PHE:HD1	3:AC:198:VAL:O	1.98	0.46
4:AD:59:ARG:NH2	4:AD:66:ARG:NH2	2.61	0.46
8:AH:103:VAL:CG1	8:AH:108:GLY:HA3	2.45	0.46
8:AH:28:ALA:O	8:AH:29:SER:HB2	2.15	0.46
9:AI:65:VAL:CG2	9:AI:66:ARG:H	2.28	0.46
1:AA:376:G:H5''	16:AP:5:ARG:CG	2.46	0.46
17:AQ:62:SER:CB	17:AQ:72:ARG:HG3	2.44	0.46
19:AS:15:LEU:CD2	19:AS:15:LEU:H	2.26	0.46
20:AT:64:ASP:C	20:AT:66:ALA:N	2.68	0.46
25:AY:147:LEU:HD23	25:AY:148:HIS:H	1.78	0.46
29:B3:26:LEU:HD21	29:B3:46:ASN:CB	2.46	0.46
34:B8:56:GLU:CA	34:B8:59:LYS:HZ1	2.28	0.46
35:BA:1689:A:O2'	35:BA:1690:A:H5'	2.14	0.46
35:BA:1827:C:H2'	35:BA:1828:G:O4'	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2087:G:C2'	35:BA:2088:G:H5'	2.46	0.46
35:BA:2400:G:H2'	35:BA:2401:U:H6	1.81	0.46
35:BA:2476:A:C2	35:BA:2477:C:C6	3.03	0.46
35:BA:1128:A:H2	35:BA:2516:G:N3	2.13	0.46
35:BA:537:C:H3'	35:BA:538:G:H8	1.80	0.46
35:BA:845:G:HO2'	35:BA:846:C:H5	1.61	0.46
35:BA:851:U:O2'	35:BA:852:G:H5'	2.16	0.46
35:BA:729:G:O5'	38:BD:208:LYS:NZ	2.48	0.46
38:BD:265:PRO:HG2	38:BD:266:SER:N	2.22	0.46
40:BF:152:GLU:OE1	40:BF:191:ARG:HD2	2.15	0.46
40:BF:192:LEU:HD21	40:BF:194:MET:CE	2.44	0.46
40:BF:9:ILE:HG12	40:BF:15:SER:N	2.30	0.46
41:BG:28:VAL:CG1	41:BG:28:VAL:O	2.62	0.46
42:BH:61:HIS:O	42:BH:65:HIS:HB3	2.15	0.46
42:BH:65:HIS:CG	42:BH:66:GLY:N	2.82	0.46
44:BN:79:PRO:HG2	44:BN:80:GLY:H	1.79	0.46
44:BN:87:LEU:O	44:BN:88:GLU:C	2.52	0.46
45:BO:104:ARG:NH1	45:BO:104:ARG:CB	2.77	0.46
45:BO:88:ASN:O	45:BO:91:LEU:N	2.44	0.46
47:BQ:34:LEU:HD11	47:BQ:129:THR:CG2	2.45	0.46
47:BQ:16:ARG:HB3	47:BQ:16:ARG:CZ	2.45	0.46
47:BQ:86:GLY:C	47:BQ:88:GLY:H	2.18	0.46
48:BR:28:LEU:HA	48:BR:34:ILE:CD1	2.46	0.46
49:BS:26:LEU:HG	49:BS:39:ILE:HD11	1.98	0.46
49:BS:95:HIS:O	49:BS:96:GLY:C	2.53	0.46
50:BT:3:ARG:O	50:BT:5:ALA:N	2.48	0.46
51:BU:59:ARG:O	51:BU:60:LEU:C	2.52	0.46
52:BV:34:GLU:HG2	52:BV:35:LEU:H	1.80	0.46
52:BV:56:SER:C	52:BV:58:VAL:H	2.19	0.46
53:BW:103:ILE:HG13	53:BW:103:ILE:O	2.14	0.46
53:BW:74:ALA:O	53:BW:75:TYR:CB	2.63	0.46
35:BA:1614:A:C6	53:BW:93:ALA:HB2	2.51	0.46
55:BY:14:LEU:CD1	55:BY:15:VAL:H	2.23	0.46
56:BZ:121:HIS:HB3	56:BZ:124:ILE:HG22	1.97	0.46
56:BZ:53:ILE:O	56:BZ:53:ILE:HG13	2.14	0.46
1:CA:106:C:O2'	1:CA:379:C:OP1	2.33	0.46
1:CA:1333:A:H2'	1:CA:1334:G:O4'	2.15	0.46
1:CA:1496:C:H2'	1:CA:1497:G:O4'	2.15	0.46
1:CA:1514:C:H2'	1:CA:1515:C:C6	2.50	0.46
1:CA:505:G:C6	1:CA:535:A:C2	3.03	0.46
1:CA:505:G:C2	1:CA:535:A:C2	3.04	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:565:U:H2'	1:CA:566:G:H8	1.80	0.46
1:CA:714:G:N2	1:CA:777:A:H1'	2.29	0.46
1:CA:874:G:H2'	1:CA:875:C:C6	2.50	0.46
1:CA:889:A:H5'	1:CA:891:U:O4'	2.16	0.46
2:CB:107:THR:HA	2:CB:110:GLN:HG3	1.98	0.46
3:CC:40:ARG:O	3:CC:41:GLY:C	2.52	0.46
4:CD:103:ASN:O	4:CD:106:TYR:N	2.48	0.46
4:CD:153:ARG:NH2	4:CD:181:MET:HG2	2.30	0.46
4:CD:199:ASN:OD1	4:CD:201:GLN:HB2	2.14	0.46
5:CE:94:ALA:HB3	5:CE:117:ASP:O	2.15	0.46
5:CE:109:ILE:HD12	5:CE:135:THR:HB	1.98	0.46
5:CE:37:ARG:HG2	5:CE:37:ARG:HH11	1.80	0.46
6:CF:16:GLN:O	6:CF:19:LEU:N	2.48	0.46
6:CF:30:LEU:HD23	6:CF:75:LEU:CD2	2.44	0.46
7:CG:25:ALA:CA	7:CG:28:ASN:HD22	2.28	0.46
8:CH:120:THR:N	8:CH:123:GLU:OE1	2.48	0.46
1:CA:1128:C:H5'	9:CI:16:ARG:NH1	2.30	0.46
9:CI:89:ASN:HB3	9:CI:92:TYR:CD1	2.51	0.46
11:CK:65:ALA:HB1	11:CK:98:LEU:CD2	2.45	0.46
12:CL:41:ARG:CG	12:CL:42:THR:H	2.28	0.46
12:CL:54:LYS:H	12:CL:54:LYS:HD2	1.79	0.46
12:CL:54:LYS:O	12:CL:70:ILE:HD13	2.15	0.46
13:CM:50:GLU:O	13:CM:54:VAL:HG23	2.14	0.46
15:CO:18:PHE:O	15:CO:19:PRO:C	2.52	0.46
17:CQ:32:TYR:O	17:CQ:34:LYS:N	2.48	0.46
1:CA:267:C:OP1	17:CQ:67:LYS:HB2	2.14	0.46
27:D1:85:LEU:HD23	27:D1:85:LEU:H	1.78	0.46
29:D3:19:GLN:C	29:D3:21:ALA:N	2.68	0.46
32:D6:11:LEU:O	32:D6:24:GLU:N	2.48	0.46
35:DA:1169:G:H1	35:DA:1180:C:H42	1.61	0.46
35:DA:814:C:H1'	35:DA:1225:G:N2	2.30	0.46
35:DA:577:G:O2'	35:DA:1254:A:OP1	2.33	0.46
35:DA:1388:G:H2'	35:DA:1389:G:H8	1.79	0.46
35:DA:1795:C:N4	35:DA:1824:G:H1	2.12	0.46
35:DA:1829:A:C6	35:DA:1830:C:C2	3.03	0.46
35:DA:1839:G:C5	35:DA:1840:G:N7	2.84	0.46
35:DA:1956:U:H2'	35:DA:1957:C:H5'	1.98	0.46
35:DA:2283:C:C5	35:DA:2389:G:H2'	2.51	0.46
35:DA:2618:G:H2'	35:DA:2619:C:H6	1.81	0.46
35:DA:2842:G:C6	35:DA:2876:G:C6	3.03	0.46
35:DA:462:C:N4	35:DA:468:G:C6	2.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:604:G:H2'	35:DA:605:C:H6	1.80	0.46
35:DA:626:U:H5''	35:DA:627:A:H5'	1.98	0.46
35:DA:76:C:H2'	35:DA:77:C:C6	2.48	0.46
35:DA:782:A:H4'	35:DA:783:A:O5'	2.15	0.46
35:DA:814:C:C5	46:DP:27:HIS:CE1	3.04	0.46
35:DA:1566:A:OP1	38:DD:211:ARG:NE	2.48	0.46
38:DD:218:ARG:HH11	38:DD:218:ARG:HG3	1.79	0.46
39:DE:119:ARG:HD2	39:DE:120:TRP:CD1	2.51	0.46
39:DE:120:TRP:CD1	39:DE:155:LYS:HB3	2.51	0.46
35:DA:2733:A:C2	39:DE:203:LYS:HA	2.50	0.46
39:DE:44:TYR:O	39:DE:45:THR:OG1	2.28	0.46
40:DF:20:LEU:HD13	40:DF:203:GLN:OE1	2.15	0.46
40:DF:34:TRP:HA	40:DF:37:VAL:HG23	1.98	0.46
41:DG:132:ASN:ND2	41:DG:157:ILE:C	2.68	0.46
41:DG:40:ASN:N	41:DG:157:ILE:HA	2.30	0.46
41:DG:6:ALA:O	41:DG:10:LYS:N	2.49	0.46
45:DO:104:ARG:HG2	45:DO:105:GLU:OE1	2.14	0.46
45:DO:23:ARG:HH11	45:DO:23:ARG:HG2	1.81	0.46
45:DO:68:GLU:CD	45:DO:78:ARG:NH1	2.67	0.46
45:DO:87:ILE:HD13	45:DO:92:GLU:C	2.36	0.46
47:DQ:48:GLU:O	47:DQ:52:VAL:HG12	2.15	0.46
47:DQ:73:PRO:HG3	47:DQ:93:TYR:CD2	2.51	0.46
35:DA:2276:G:OP2	47:DQ:84:GLY:N	2.47	0.46
50:DT:30:VAL:O	50:DT:31:SER:HB3	2.15	0.46
50:DT:96:ARG:HG2	50:DT:96:ARG:NH1	2.25	0.46
50:DT:96:ARG:CG	50:DT:96:ARG:NH1	2.75	0.46
52:DV:56:SER:C	52:DV:58:VAL:H	2.18	0.46
54:DX:26:TYR:CD1	54:DX:26:TYR:N	2.83	0.46
56:DZ:146:ILE:HG22	56:DZ:174:VAL:CG1	2.42	0.46
56:DZ:150:LEU:N	56:DZ:150:LEU:HD22	2.27	0.46
1:AA:1210:C:H4'	1:AA:1214:C:C4	2.50	0.46
1:AA:1332:A:C2	1:AA:1333:A:C5	3.03	0.46
1:AA:1376:U:O2'	1:AA:1377:A:H5'	2.16	0.46
1:AA:1430:C:H2'	1:AA:1431:C:H6	1.80	0.46
1:AA:152:A:N6	1:AA:170:U:N3	2.64	0.46
1:AA:189(I):G:H2'	1:AA:189(J):G:C8	2.50	0.46
1:AA:261:U:O2	1:AA:263:A:C8	2.68	0.46
1:AA:325:A:N6	1:AA:326:G:C6	2.83	0.46
1:AA:514:C:O2'	1:AA:515:G:H5'	2.16	0.46
1:AA:652:U:H1'	1:AA:653:A:C2	2.47	0.46
1:AA:786:G:C2	1:AA:797:C:O2	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:12:GLU:C	2:AB:14:GLY:H	2.17	0.46
2:AB:55:PHE:HA	2:AB:58:ILE:CB	2.42	0.46
4:AD:23:GLY:O	4:AD:27:TYR:CD1	2.68	0.46
6:AF:88:VAL:HG12	6:AF:89:MET:N	2.31	0.46
7:AG:50:ILE:CB	7:AG:58:PRO:HD3	2.40	0.46
9:AI:56:LEU:HD23	9:AI:56:LEU:O	2.15	0.46
9:AI:28:VAL:HG22	9:AI:63:ILE:O	2.15	0.46
12:AL:60:LEU:HD23	12:AL:64:TYR:C	2.36	0.46
16:AP:77:ALA:HB3	16:AP:79:VAL:HG23	1.98	0.46
16:AP:80:PHE:O	16:AP:81:ARG:C	2.53	0.46
20:AT:25:ARG:HB2	20:AT:25:ARG:CZ	2.45	0.46
22:AV:40:C:H2'	22:AV:41:C:C6	2.49	0.46
25:AY:67:VAL:HG12	25:AY:100:TYR:CD1	2.49	0.46
25:AY:129:ILE:HG21	25:AY:169:ILE:HG12	1.98	0.46
25:AY:162:GLN:O	25:AY:166:ASP:OD2	2.33	0.46
27:B1:11:ARG:HG3	27:B1:61:ARG:O	2.14	0.46
28:B2:22:GLU:O	28:B2:25:VAL:CG1	2.64	0.46
35:BA:1186:G:H2'	35:BA:1187:G:C5'	2.45	0.46
35:BA:1210:A:H1'	35:BA:1212:G:N3	2.31	0.46
35:BA:1472:A:O2'	35:BA:1473:G:H5'	2.15	0.46
35:BA:1464:C:H1'	35:BA:1528(A):A:N3	2.31	0.46
35:BA:1638:C:H2'	35:BA:1639:U:C6	2.50	0.46
35:BA:1644:C:C2'	35:BA:1645:G:H5'	2.45	0.46
35:BA:1766:U:H2'	35:BA:1767:C:H6	1.80	0.46
35:BA:2014:A:C2	35:BA:2015:A:N1	2.83	0.46
35:BA:2039:C:H2'	35:BA:2040:C:C6	2.50	0.46
35:BA:2240:C:O2'	35:BA:2241:A:H5'	2.15	0.46
35:BA:483:A:H1'	55:BY:47:LYS:CG	2.24	0.46
35:BA:559:G:H2'	35:BA:560:C:H6	1.80	0.46
35:BA:804:A:H2'	35:BA:806:C:C4	2.51	0.46
35:BA:814:C:H2'	35:BA:815:C:H6	1.80	0.46
36:BB:28:C:H2'	36:BB:29:A:O4'	2.16	0.46
35:BA:1971:A:C4	38:BD:241:PRO:HB3	2.50	0.46
39:BE:199:ARG:HB3	39:BE:200:GLU:OE2	2.14	0.46
35:BA:2733:A:C2	39:BE:203:LYS:HA	2.49	0.46
40:BF:21:ALA:C	40:BF:23:ASP:N	2.68	0.46
41:BG:131:TYR:CE1	41:BG:133:LEU:HB3	2.51	0.46
43:BI:71:ILE:HG13	43:BI:72:LEU:HD22	1.98	0.46
45:BO:37:ASP:HB2	45:BO:62:VAL:HG23	1.96	0.46
45:BO:68:GLU:CD	45:BO:78:ARG:NH1	2.66	0.46
46:BP:96:THR:HB	46:BP:126:VAL:HB	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1202:C:O2	46:BP:7:ARG:NH2	2.49	0.46
48:BR:32:GLY:O	48:BR:116:LEU:HB2	2.15	0.46
48:BR:98:LEU:HB2	48:BR:113:LEU:CD2	2.30	0.46
49:BS:106:ARG:HB3	49:BS:106:ARG:HE	1.52	0.46
51:BU:78:THR:C	51:BU:80:ILE:H	2.18	0.46
28:B2:26:ARG:HH21	54:BX:6:ASP:HA	1.80	0.46
56:BZ:126:VAL:C	56:BZ:164:ALA:HB3	2.35	0.46
1:CA:1011:G:C2'	1:CA:1012:U:H5'	2.45	0.46
1:CA:1217:C:H2'	1:CA:1218:C:C6	2.51	0.46
1:CA:1523:G:C6	1:CA:1524:C:C4	3.04	0.46
1:CA:237:C:C5'	17:CQ:25:ARG:NH1	2.78	0.46
1:CA:302:G:N3	1:CA:556:C:H4'	2.30	0.46
1:CA:839:U:O2	1:CA:839:U:O4'	2.34	0.46
2:CB:15:VAL:HG23	2:CB:209:ARG:HE	1.80	0.46
2:CB:88:ALA:CB	2:CB:223:ILE:HD11	2.44	0.46
4:CD:3:ARG:O	4:CD:4:TYR:C	2.53	0.46
4:CD:62:GLN:HB3	4:CD:66:ARG:HH12	1.79	0.46
6:CF:12:PRO:HG2	6:CF:13:ASN:H	1.80	0.46
6:CF:21:LEU:O	6:CF:23:LYS:N	2.48	0.46
7:CG:146:GLU:CA	7:CG:149:ARG:HB2	2.43	0.46
7:CG:42:ILE:HG23	7:CG:117:ALA:CA	2.46	0.46
1:CA:1349:A:OP2	9:CI:118:LYS:NZ	2.48	0.46
9:CI:28:VAL:CG2	9:CI:63:ILE:HB	2.45	0.46
15:CO:39:LEU:HD12	15:CO:56:LEU:HD12	1.97	0.46
18:CR:76:LEU:HD22	18:CR:76:LEU:N	2.30	0.46
23:CW:4:G:C6	23:CW:71:G:C6	3.04	0.46
25:CY:109:GLU:HA	25:CY:112:LYS:HB3	1.97	0.46
27:D1:51:VAL:O	27:D1:59:THR:HA	2.15	0.46
28:D2:15:LYS:C	28:D2:18:PRO:HD2	2.35	0.46
35:DA:1131:G:C2	35:DA:1132:A:C5	3.03	0.46
35:DA:1274:A:C2	35:DA:1302:A:C2	3.04	0.46
35:DA:1496:A:H8	35:DA:1577:C:O2'	1.98	0.46
35:DA:1784:A:H4'	35:DA:1785:A:O5'	2.16	0.46
35:DA:2299:G:C2	35:DA:2318:G:C8	3.03	0.46
33:D7:34:ARG:NH1	35:DA:466:A:OP1	2.48	0.46
35:DA:481:G:H2'	35:DA:507:A:N1	2.31	0.46
35:DA:538:G:H2'	35:DA:538:G:N3	2.30	0.46
35:DA:601:C:O2'	35:DA:605:C:H5''	2.16	0.46
35:DA:621:A:H2'	35:DA:622:G:C5'	2.44	0.46
36:DB:57:A:C1'	41:DG:30:GLU:HB3	2.45	0.46
38:DD:165:ILE:C	38:DD:166:GLN:NE2	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:25:THR:O	38:DD:26:LYS:C	2.53	0.46
38:DD:267:SER:O	38:DD:269:PHE:HD1	1.98	0.46
38:DD:63:ARG:CZ	38:DD:86:PRO:HD2	2.45	0.46
40:DF:22:ALA:HA	40:DF:26:ALA:HB2	1.97	0.46
41:DG:39:ILE:HG13	41:DG:157:ILE:CG2	2.36	0.46
41:DG:98:ARG:O	41:DG:101:ILE:CG2	2.63	0.46
42:DH:85:LYS:HZ3	42:DH:145:ALA:HA	1.80	0.46
44:DN:128:HIS:C	44:DN:130:HIS:HD1	2.18	0.46
44:DN:43:THR:O	44:DN:46:VAL:HG12	2.14	0.46
45:DO:87:ILE:HG13	45:DO:91:LEU:HD13	1.96	0.46
35:DA:626:U:C2	46:DP:105:LEU:HG	2.50	0.46
46:DP:48:PRO:HG2	46:DP:49:ARG:N	2.30	0.46
46:DP:65:ARG:HG2	46:DP:65:ARG:O	2.15	0.46
47:DQ:113:GLN:O	47:DQ:116:GLU:HB3	2.15	0.46
48:DR:101:ALA:HB1	53:DW:38:TYR:HE1	1.80	0.46
48:DR:9:LYS:HZ3	48:DR:42:LYS:HB3	1.79	0.46
50:DT:28:VAL:O	50:DT:29:ARG:CD	2.64	0.46
50:DT:88:ILE:HG22	50:DT:89:VAL:H	1.78	0.46
51:DU:98:LEU:HD22	52:DV:2:PHE:HZ	1.78	0.46
52:DV:75:PHE:HD1	52:DV:87:HIS:O	1.98	0.46
54:DX:35:THR:O	54:DX:36:LYS:C	2.54	0.46
54:DX:64:LYS:CG	54:DX:65:ARG:H	2.28	0.46
54:DX:63:LYS:O	54:DX:68:ARG:HA	2.16	0.46
56:DZ:137:ILE:HD13	56:DZ:155:LEU:HD13	1.98	0.46
1:AA:112:G:H5'	1:AA:389:A:H4'	1.97	0.46
1:AA:1153:C:H2'	1:AA:1154:G:O5'	2.15	0.46
1:AA:286:G:H2'	1:AA:287:U:H6	1.80	0.46
1:AA:712:A:O2'	1:AA:713:G:H5'	2.15	0.46
1:AA:807:A:H2'	1:AA:808:C:C6	2.50	0.46
2:AB:36:ARG:O	2:AB:37:ASN:C	2.53	0.46
3:AC:105:GLU:HG2	3:AC:106:VAL:N	2.19	0.46
5:AE:128:PRO:O	5:AE:131:ILE:HG12	2.15	0.46
5:AE:144:THR:O	5:AE:145:LYS:C	2.54	0.46
7:AG:145:ALA:O	7:AG:147:ALA:N	2.36	0.46
9:AI:17:VAL:HG21	9:AI:81:ILE:HD13	1.97	0.46
9:AI:37:PHE:HB3	9:AI:43:ALA:HB2	1.97	0.46
12:AL:28:LYS:C	12:AL:30:ALA:N	2.69	0.46
12:AL:70:ILE:H	12:AL:70:ILE:HD12	1.78	0.46
1:AA:658:G:O4'	15:AO:22:THR:HB	2.14	0.46
15:AO:29:VAL:O	15:AO:30:ALA:C	2.53	0.46
15:AO:37:ASN:HD22	15:AO:37:ASN:N	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AO:47:LYS:HD3	15:AO:47:LYS:H	1.80	0.46
7:AG:86:GLN:CD	23:AW:32:G:H21	2.18	0.46
27:B1:48:LYS:CG	27:B1:49:VAL:N	2.76	0.46
29:B3:16:PRO:HB2	29:B3:18:ASP:OD1	2.16	0.46
33:B7:18:PHE:O	33:B7:19:ARG:C	2.54	0.46
35:BA:1169:G:H1	35:BA:1180:C:H42	1.64	0.46
35:BA:1338:G:N3	35:BA:1338:G:H2'	2.29	0.46
35:BA:1433:U:H3	35:BA:1560:G:H1	1.62	0.46
35:BA:1598:C:H5'	54:BX:37:THR:HB	1.96	0.46
35:BA:1349:A:N6	35:BA:1598:C:N4	2.62	0.46
35:BA:1789:A:H2'	35:BA:1790:C:C6	2.48	0.46
35:BA:2033:A:O2'	35:BA:2034:U:P	2.73	0.46
35:BA:2208:A:H1'	35:BA:2219:G:C5	2.50	0.46
35:BA:237:C:O2'	35:BA:238:C:H5'	2.15	0.46
35:BA:2866:U:C6	35:BA:2868:A:C1'	2.98	0.46
35:BA:315:G:H2'	35:BA:316:C:C6	2.50	0.46
35:BA:483:A:N7	35:BA:497:A:H2	2.13	0.46
35:BA:558:G:H2'	35:BA:559:G:C8	2.44	0.46
35:BA:645:C:C2'	35:BA:645:C:O2	2.64	0.46
35:BA:869:G:H2'	35:BA:870:A:C8	2.49	0.46
35:BA:948:G:H2'	35:BA:949:C:C6	2.51	0.46
36:BB:79:C:H2'	36:BB:80:U:O4'	2.15	0.46
40:BF:160:ASN:HD21	40:BF:162:LEU:CB	2.29	0.46
40:BF:170:LEU:HD21	40:BF:172:TRP:CZ2	2.50	0.46
44:BN:94:HIS:N	44:BN:94:HIS:ND1	2.64	0.46
45:BO:64:ARG:HD3	45:BO:101:PRO:C	2.35	0.46
46:BP:101:VAL:HA	46:BP:107:LYS:H	1.79	0.46
47:BQ:35:VAL:CG2	47:BQ:101:ARG:O	2.63	0.46
49:BS:97:ARG:O	49:BS:98:VAL:C	2.53	0.46
44:BN:40:PRO:CA	51:BU:64:ARG:HH22	2.28	0.46
51:BU:65:ILE:H	51:BU:65:ILE:CD1	2.28	0.46
51:BU:92:ARG:HG2	51:BU:92:ARG:O	2.15	0.46
52:BV:38:LEU:O	52:BV:53:GLU:O	2.33	0.46
53:BW:48:ALA:O	53:BW:49:LYS:C	2.54	0.46
53:BW:65:LEU:HD23	53:BW:65:LEU:C	2.35	0.46
28:B2:51:ARG:NH2	54:BX:6:ASP:OD2	2.49	0.46
56:BZ:19:ARG:HG3	56:BZ:20:ARG:N	2.30	0.46
1:CA:1133:G:H2'	1:CA:1134:G:C8	2.50	0.46
1:CA:264:U:O2'	17:CQ:64:PRO:HB2	2.16	0.46
1:CA:39:G:H2'	1:CA:40:C:H6	1.79	0.46
1:CA:751:U:O2'	1:CA:752:G:H5'	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:68:ILE:HG22	2:CB:70:PHE:CD1	2.50	0.46
3:CC:11:ARG:O	3:CC:14:ILE:O	2.34	0.46
4:CD:147:ALA:HA	4:CD:182:LYS:HB3	1.97	0.46
4:CD:23:GLY:O	4:CD:27:TYR:CD1	2.69	0.46
4:CD:65:ARG:NH1	4:CD:72:GLU:HB2	2.31	0.46
7:CG:140:ASP:HA	7:CG:143:ARG:CZ	2.46	0.46
8:CH:86:ILE:CG1	8:CH:135:CYS:HA	2.41	0.46
9:CI:71:SER:O	9:CI:74:ILE:HB	2.15	0.46
10:CJ:31:GLY:HA3	10:CJ:78:ASN:ND2	2.29	0.46
10:CJ:23:ILE:HG23	10:CJ:85:LEU:HD13	1.97	0.46
10:CJ:96:ILE:HD13	10:CJ:96:ILE:N	2.30	0.46
11:CK:15:ALA:HA	11:CK:76:GLY:O	2.15	0.46
12:CL:18:VAL:HG23	12:CL:18:VAL:O	2.16	0.46
17:CQ:67:LYS:C	17:CQ:69:LYS:H	2.19	0.46
19:CS:80:TYR:O	19:CS:81:ARG:O	2.34	0.46
23:CW:70:C:H2'	23:CW:71:G:H8	1.80	0.46
25:CY:7:TYR:CE1	25:CY:160:GLU:HG2	2.50	0.46
35:DA:1149:G:H2'	35:DA:1150:C:C6	2.51	0.46
35:DA:1272:A:N3	35:DA:1618:A:C5	2.83	0.46
35:DA:1281:G:H2'	35:DA:1282:U:H6	1.79	0.46
35:DA:1630:G:C5	35:DA:1631:C:C5	3.04	0.46
35:DA:1668:A:C8	35:DA:1674:G:C6	3.04	0.46
35:DA:2007:C:H6	35:DA:2007:C:O5'	1.97	0.46
35:DA:200:U:C2'	35:DA:201:C:H5'	2.44	0.46
35:DA:2023:G:H2'	35:DA:2024:G:H8	1.81	0.46
35:DA:2030:A:H4'	35:DA:2031:A:C8	2.47	0.46
35:DA:2039:C:C2	35:DA:2040:C:C5	3.04	0.46
35:DA:2124:G:H2'	35:DA:2125:G:H5'	1.97	0.46
35:DA:2171:A:C8	35:DA:2172:U:C5	3.04	0.46
35:DA:2362:G:O2'	35:DA:2363:C:H5'	2.15	0.46
35:DA:2473:U:O2	35:DA:2473:U:C2'	2.62	0.46
35:DA:2577:A:H5''	35:DA:2578:G:C5'	2.40	0.46
35:DA:2643:G:C2'	35:DA:2644:G:H5'	2.45	0.46
35:DA:322:A:OP1	40:DF:168:ARG:HD3	2.16	0.46
35:DA:419:C:C2	35:DA:420:C:C6	3.04	0.46
35:DA:448:U:H4'	35:DA:449:A:OP2	2.15	0.46
38:DD:227:ASN:O	38:DD:229:VAL:N	2.49	0.46
39:DE:120:TRP:O	39:DE:121:ASN:C	2.54	0.46
39:DE:93:VAL:C	39:DE:95:ILE:N	2.68	0.46
40:DF:114:VAL:O	40:DF:115:ALA:C	2.52	0.46
40:DF:160:ASN:HD21	40:DF:162:LEU:CB	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DF:117:ARG:NH2	40:DF:187:VAL:HA	2.30	0.46
40:DF:2:LYS:CG	40:DF:25:PRO:HB2	2.22	0.46
43:DI:98:ALA:HB1	43:DI:109:ILE:HD13	1.96	0.46
44:DN:9:VAL:HG21	44:DN:39:ARG:HH21	1.80	0.46
46:DP:101:VAL:O	46:DP:103:ALA:N	2.49	0.46
46:DP:38:GLN:CG	46:DP:39:LYS:N	2.71	0.46
47:DQ:82:ARG:HH11	47:DQ:82:ARG:CG	2.25	0.46
49:DS:92:TYR:HD2	49:DS:97:ARG:NH1	2.12	0.46
51:DU:36:ARG:HG2	51:DU:40:PHE:CE2	2.50	0.46
51:DU:76:TYR:CZ	51:DU:80:ILE:HG13	2.50	0.46
51:DU:83:LEU:HG	51:DU:88:ILE:HG12	1.96	0.46
52:DV:75:PHE:HE1	52:DV:89:GLN:HB2	1.79	0.46
53:DW:20:VAL:HG23	53:DW:21:VAL:H	1.80	0.46
54:DX:12:VAL:HG13	54:DX:17:ALA:HB2	1.98	0.46
1:AA:1039:C:H2'	1:AA:1040:U:C6	2.51	0.46
1:AA:1120:G:H1	1:AA:1153:C:H42	1.63	0.46
1:AA:1251:A:H2'	1:AA:1252:A:C8	2.50	0.46
1:AA:1255:G:O2'	1:AA:1258:G:H1'	2.16	0.46
1:AA:1423:G:N1	1:AA:1478:C:N3	2.64	0.46
1:AA:182:U:H2'	1:AA:183:G:O4'	2.15	0.46
1:AA:193:C:H2'	1:AA:194:C:C6	2.50	0.46
1:AA:35:G:H2'	1:AA:36:C:H6	1.77	0.46
1:AA:799:G:C2'	1:AA:800:G:H5'	2.46	0.46
1:AA:929:G:C6	1:AA:930:C:C4	3.04	0.46
2:AB:111:ARG:HH11	2:AB:111:ARG:CG	2.22	0.46
2:AB:20:GLU:OE1	2:AB:20:GLU:HA	2.15	0.46
2:AB:80:ILE:HG12	2:AB:211:ILE:CG2	2.46	0.46
2:AB:55:PHE:O	2:AB:59:GLU:N	2.41	0.46
3:AC:92:ALA:C	3:AC:94:LEU:H	2.17	0.46
4:AD:152:SER:C	4:AD:154:ASN:H	2.19	0.46
4:AD:178:VAL:C	4:AD:180:GLY:H	2.18	0.46
5:AE:136:MET:HB3	5:AE:140:ARG:HH21	1.79	0.46
7:AG:81:GLY:C	7:AG:83:ALA:H	2.17	0.46
11:AK:73:MET:SD	11:AK:103:LEU:CD1	3.03	0.46
1:AA:568:G:O6	12:AL:5:PRO:HD3	2.16	0.46
15:AO:16:ALA:HB1	15:AO:21:ASP:HB3	1.97	0.46
18:AR:43:PHE:CA	18:AR:51:LEU:HD12	2.37	0.46
19:AS:33:THR:CG2	19:AS:51:VAL:HG22	2.46	0.46
20:AT:13:LEU:HD13	20:AT:14:LYS:N	2.31	0.46
20:AT:14:LYS:HE3	20:AT:18:GLN:HE21	1.80	0.46
1:AA:263:A:OP2	20:AT:79:ARG:NH1	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:790:A:OP1	22:AV:37:A:O2'	2.31	0.46
25:AY:53:ASN:HD22	25:AY:53:ASN:H	1.63	0.46
25:AY:6:LEU:O	25:AY:6:LEU:HD23	2.15	0.46
31:B5:49:CYS:O	31:B5:57:VAL:HG22	2.15	0.46
35:BA:1221(A):C:C2'	35:BA:1222:C:H5'	2.45	0.46
35:BA:1388:G:H2'	35:BA:1389:G:H8	1.80	0.46
35:BA:1531:C:H3'	35:BA:1532:C:C5'	2.46	0.46
35:BA:1817:G:C2'	35:BA:1818:U:H5'	2.46	0.46
35:BA:1935:G:H1'	35:BA:1964:G:H21	1.78	0.46
35:BA:197:A:C2	35:BA:198:C:H1'	2.51	0.46
35:BA:2376:A:H2'	35:BA:2377:A:O4'	2.15	0.46
35:BA:2447:G:O6	35:BA:2504:U:O4	2.34	0.46
35:BA:2643:G:C2'	35:BA:2644:G:H5'	2.45	0.46
35:BA:526:A:OP1	35:BA:527:C:OP1	2.33	0.46
35:BA:587:C:O2'	35:BA:588:U:OP2	2.26	0.46
35:BA:680:G:O2'	35:BA:681:G:H5'	2.16	0.46
35:BA:782:A:H4'	35:BA:783:A:O5'	2.16	0.46
35:BA:874:G:H2'	35:BA:875:G:C8	2.50	0.46
38:BD:27:THR:CG2	38:BD:28:GLU:N	2.62	0.46
39:BE:108:SER:O	39:BE:162:ALA:HA	2.15	0.46
39:BE:44:TYR:O	39:BE:45:THR:OG1	2.28	0.46
39:BE:73:GLU:OE2	39:BE:73:GLU:N	2.34	0.46
39:BE:78:LEU:O	39:BE:79:ARG:HD2	2.15	0.46
40:BF:126:VAL:CG1	40:BF:193:VAL:HG13	2.46	0.46
41:BG:138:GLN:HE21	41:BG:152:LEU:HD23	1.81	0.46
41:BG:36:LYS:HE3	41:BG:160:VAL:CG2	2.46	0.46
42:BH:162:ILE:C	42:BH:163:TYR:CD1	2.89	0.46
42:BH:97:ARG:O	42:BH:98:LEU:HB2	2.15	0.46
44:BN:31:ALA:O	44:BN:34:LEU:N	2.49	0.46
45:BO:104:ARG:C	45:BO:106:LEU:H	2.17	0.46
45:BO:113:LYS:O	45:BO:114:ILE:C	2.54	0.46
45:BO:112:MET:O	45:BO:115:VAL:HG23	2.16	0.46
46:BP:14:LYS:O	46:BP:15:ARG:HB2	2.16	0.46
47:BQ:74:TYR:HB3	47:BQ:91:GLU:OE2	2.15	0.46
36:BB:7:G:H21	49:BS:38:GLN:HE22	1.63	0.46
51:BU:111:GLU:O	51:BU:113:ALA:N	2.49	0.46
51:BU:26:GLY:C	51:BU:28:ARG:H	2.18	0.46
52:BV:17:GLY:HA2	52:BV:98:GLU:O	2.14	0.46
1:CA:1065:U:H4'	1:CA:1066:C:O5'	2.16	0.46
1:CA:267:C:P	17:CQ:67:LYS:HB2	2.56	0.46
1:CA:487:A:C6	1:CA:488:C:O2	2.69	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:528:C:C5	1:CA:529:G:N7	2.84	0.46
1:CA:730:G:C5	1:CA:731:G:H1'	2.51	0.46
1:CA:763:G:H2'	1:CA:764:C:H6	1.80	0.46
1:CA:818:G:C3'	1:CA:819:A:C5'	2.91	0.46
2:CB:99:GLY:O	2:CB:100:GLY:C	2.52	0.46
2:CB:92:TYR:CE2	2:CB:151:GLY:HA3	2.51	0.46
3:CC:84:ILE:HG13	3:CC:101:LEU:HD13	1.96	0.46
3:CC:155:GLY:HA3	3:CC:163:ALA:CB	2.42	0.46
7:CG:107:ALA:O	7:CG:110:GLN:HB2	2.15	0.46
8:CH:5:PRO:C	8:CH:8:ASP:HB3	2.36	0.46
12:CL:89:ARG:CB	12:CL:89:ARG:HH11	2.28	0.46
12:CL:89:ARG:NH1	12:CL:90:VAL:N	2.63	0.46
14:CN:24:CYS:SG	14:CN:39:LEU:HA	2.56	0.46
19:CS:67:VAL:HG12	19:CS:68:GLY:N	2.31	0.46
20:CT:27:LYS:O	20:CT:30:LYS:HB3	2.15	0.46
24:CX:17:U:H2'	24:CX:18:C:H6	1.80	0.46
25:CY:21:LEU:O	25:CY:22:GLU:C	2.54	0.46
25:CY:3:LEU:HA	25:CY:6:LEU:CB	2.29	0.46
26:D0:84:LEU:C	26:D0:84:LEU:HD12	2.35	0.46
28:D2:44:LEU:HA	28:D2:44:LEU:HD12	1.57	0.46
35:DA:1231:G:H2'	35:DA:1232:G:H8	1.81	0.46
35:DA:1266:G:N2	35:DA:2012:G:C4	2.83	0.46
35:DA:1271:G:C2	35:DA:1617:C:H4'	2.50	0.46
35:DA:1272:A:C2	35:DA:1618:A:C5	3.03	0.46
35:DA:1399:C:H2'	35:DA:1400:G:C8	2.49	0.46
35:DA:1675:C:H2'	35:DA:1676:A:O4'	2.15	0.46
35:DA:1889:A:H1'	35:DA:2087:G:O4'	2.15	0.46
35:DA:2014:A:C2	35:DA:2015:A:C2	3.03	0.46
35:DA:2791:C:H6	35:DA:2793:G:O6	1.97	0.46
35:DA:2792:G:O6	35:DA:2804:C:N3	2.48	0.46
35:DA:572:A:H2'	35:DA:573:G:O4'	2.16	0.46
35:DA:923:C:H2'	35:DA:924:C:C6	2.51	0.46
35:DA:976:C:H2'	35:DA:977:G:H8	1.79	0.46
38:DD:109:ASP:N	38:DD:195:ALA:O	2.30	0.46
38:DD:267:SER:O	38:DD:269:PHE:CD1	2.69	0.46
40:DF:117:ARG:NH2	40:DF:189:THR:O	2.48	0.46
40:DF:21:ALA:C	40:DF:23:ASP:H	2.18	0.46
40:DF:65:TRP:CZ3	40:DF:72:ARG:HB3	2.51	0.46
42:DH:13:LYS:C	42:DH:15:VAL:N	2.65	0.46
43:DI:88:ILE:CD1	43:DI:123:LEU:HD12	2.41	0.46
43:DI:39:ALA:O	43:DI:40:THR:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:122:VAL:CG1	44:DN:123:TYR:N	2.78	0.46
44:DN:67:LEU:HB3	44:DN:88:GLU:CD	2.36	0.46
45:DO:64:ARG:CB	45:DO:64:ARG:HH11	2.29	0.46
45:DO:61:VAL:CG1	45:DO:85:VAL:HB	2.45	0.46
50:DT:65:LYS:HE3	50:DT:66:VAL:H	1.80	0.46
50:DT:83:ILE:CG1	50:DT:84:GLN:HG2	2.44	0.46
53:DW:107:LEU:HA	53:DW:107:LEU:HD12	1.73	0.46
53:DW:69:LEU:N	53:DW:69:LEU:HD12	2.31	0.46
54:DX:37:THR:C	54:DX:39:ILE:H	2.19	0.46
28:D2:26:ARG:HH21	54:DX:7:VAL:H	1.60	0.46
56:DZ:76:LEU:H	56:DZ:76:LEU:HD23	1.80	0.46
56:DZ:8:TYR:HA	56:DZ:62:PRO:CG	2.45	0.46
1:AA:1002:G:H2'	1:AA:1003:G:O4'	2.15	0.46
1:AA:1161:C:H2'	1:AA:1162:C:H6	1.81	0.46
1:AA:1049:U:H1'	1:AA:1201:A:C8	2.51	0.46
1:AA:1226:C:H5'	19:AS:80:TYR:HE2	1.80	0.46
1:AA:1277:C:H2'	1:AA:1278:U:H5'	1.97	0.46
1:AA:1312:G:H1	1:AA:1325:C:H42	1.61	0.46
1:AA:236:G:H2'	1:AA:237:C:H6	1.80	0.46
1:AA:450:G:OP1	16:AP:43:LYS:NZ	2.49	0.46
1:AA:598:U:H4'	8:AH:94:TYR:CD1	2.51	0.46
1:AA:762:C:H2'	1:AA:763:G:H8	1.81	0.46
1:AA:769:G:C4	1:AA:770:C:C5	3.04	0.46
2:AB:77:ALA:C	2:AB:80:ILE:HG23	2.36	0.46
4:AD:3:ARG:HD3	4:AD:3:ARG:O	2.15	0.46
6:AF:21:LEU:O	6:AF:23:LYS:N	2.48	0.46
6:AF:37:VAL:HG13	6:AF:65:VAL:HG12	1.96	0.46
7:AG:99:LEU:O	7:AG:100:ALA:C	2.54	0.46
1:AA:376:G:C5'	16:AP:5:ARG:HD2	2.44	0.46
17:AQ:45:HIS:NE2	17:AQ:47:PRO:HB3	2.30	0.46
17:AQ:64:PRO:C	17:AQ:65:ILE:HD12	2.36	0.46
19:AS:78:ARG:H	19:AS:78:ARG:CD	2.28	0.46
19:AS:80:TYR:O	19:AS:81:ARG:O	2.34	0.46
25:AY:14:MET:HB3	25:AY:168:PHE:CD2	2.51	0.46
25:AY:30:THR:OG1	25:AY:31:GLY:N	2.49	0.46
25:AY:38:LEU:O	25:AY:41:LEU:HB2	2.16	0.46
27:B1:11:ARG:C	27:B1:13:ILE:H	2.19	0.46
33:B7:34:ARG:HE	33:B7:39:ARG:CD	2.29	0.46
35:BA:1022:G:O2'	35:BA:1023:U:P	2.73	0.46
35:BA:1110:G:OP1	35:BA:1110:G:H4'	2.16	0.46
35:BA:1459:G:H5''	35:BA:1460:A:OP2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1706:U:O2'	35:BA:1707:G:H5'	2.15	0.46
35:BA:1708:C:O2'	35:BA:1709:U:H5'	2.15	0.46
35:BA:1806:C:N4	35:BA:1812:A:N6	2.63	0.46
35:BA:1848:A:H2'	35:BA:1849:G:C8	2.50	0.46
35:BA:571:A:H5''	35:BA:2030:A:H62	1.79	0.46
35:BA:2124:G:H2'	35:BA:2125:G:H5'	1.97	0.46
35:BA:2322:A:O2'	35:BA:2323:G:H5'	2.16	0.46
35:BA:2517:C:H2'	35:BA:2542:A:H2	1.78	0.46
35:BA:2681:C:C4	35:BA:2724:C:C5	3.04	0.46
35:BA:2840:C:H2'	35:BA:2841:C:C6	2.50	0.46
35:BA:2884:U:H2'	35:BA:2885:C:C5'	2.44	0.46
35:BA:365:C:H6	35:BA:365:C:H5'	1.79	0.46
35:BA:469:G:C2'	35:BA:470:A:H5''	2.46	0.46
35:BA:532:A:H4'	35:BA:533:G:O4'	2.16	0.46
35:BA:707:G:H3'	35:BA:708:C:H6	1.80	0.46
36:BB:115:G:H2'	36:BB:116:G:H8	1.81	0.46
38:BD:145:VAL:CG1	38:BD:146:GLU:N	2.78	0.46
38:BD:208:LYS:O	38:BD:210:GLY:O	2.34	0.46
38:BD:223:GLY:C	38:BD:225:ALA:H	2.14	0.46
38:BD:265:PRO:O	38:BD:266:SER:C	2.53	0.46
40:BF:108:LYS:O	40:BF:112:MET:SD	2.74	0.46
40:BF:123:LEU:HD12	40:BF:124:LEU:N	2.29	0.46
40:BF:125:LEU:HD11	40:BF:199:TRP:CD1	2.51	0.46
41:BG:76:SER:HB3	41:BG:84:LYS:CD	2.45	0.46
43:BI:90:GLY:O	43:BI:91:SER:O	2.34	0.46
45:BO:25:LEU:HD13	45:BO:38:VAL:HG11	1.97	0.46
49:BS:101:LEU:HD21	49:BS:103:GLU:HG3	1.97	0.46
54:BX:24:GLY:HA2	54:BX:80:ILE:HG13	1.96	0.46
1:CA:1046:A:H2'	1:CA:1046:A:N3	2.30	0.46
1:CA:1051:C:H2'	1:CA:1052:U:H6	1.80	0.46
1:CA:1366:C:H2'	1:CA:1367:C:H6	1.75	0.46
1:CA:1508:G:C4	1:CA:1509:C:C5	3.03	0.46
1:CA:648:A:H2'	1:CA:649:G:C8	2.47	0.46
1:CA:577:G:H1'	1:CA:816:A:N3	2.29	0.46
1:CA:826:C:H2'	1:CA:827:U:H6	1.79	0.46
1:CA:835:U:O2'	1:CA:836:G:H5'	2.15	0.46
1:CA:957:U:H1'	1:CA:960:U:C4	2.50	0.46
1:CA:959:A:H3'	1:CA:960:U:H5''	1.97	0.46
2:CB:46:LYS:O	2:CB:47:THR:C	2.53	0.46
2:CB:84:GLU:HB3	2:CB:219:VAL:CG2	2.32	0.46
1:CA:438:G:H4'	4:CD:123:HIS:ND1	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:148:VAL:HG23	4:CD:181:MET:O	2.16	0.46
8:CH:64:LYS:O	8:CH:79:VAL:CG2	2.63	0.46
9:CI:5:TYR:OH	9:CI:7:THR:HG23	2.16	0.46
10:CJ:57:LYS:HD2	10:CJ:60:ARG:NH2	2.30	0.46
13:CM:69:GLU:HB2	13:CM:72:ALA:HB3	1.98	0.46
10:CJ:65:LEU:HD12	14:CN:55:GLY:HA3	1.98	0.46
15:CO:10:LYS:HG3	15:CO:11:VAL:N	2.30	0.46
15:CO:54:ARG:NH1	15:CO:54:ARG:HG2	2.30	0.46
19:CS:17:GLU:C	19:CS:19:VAL:H	2.18	0.46
25:CY:18:LEU:C	25:CY:20:VAL:H	2.19	0.46
35:DA:1198:U:H2'	35:DA:1199:U:H6	1.80	0.46
35:DA:1396:U:O2	35:DA:1396:U:C2'	2.64	0.46
35:DA:1516:C:H2'	35:DA:1517:G:C8	2.51	0.46
35:DA:1567:A:H3'	38:DD:86:PRO:HG3	1.97	0.46
35:DA:1689:A:O2'	35:DA:1690:A:H5'	2.15	0.46
35:DA:1692:U:O2'	35:DA:1693:U:H2'	2.15	0.46
35:DA:1695:G:H2'	35:DA:1696:G:C4'	2.45	0.46
27:D1:25:LYS:NZ	35:DA:2079:U:OP1	2.49	0.46
35:DA:2219:G:C2'	35:DA:2220:G:H5'	2.45	0.46
35:DA:2291:U:H2'	35:DA:2292:C:C6	2.50	0.46
35:DA:2377:A:H2'	35:DA:2378:A:C8	2.51	0.46
35:DA:2285:C:H42	35:DA:2383:G:H1	1.64	0.46
35:DA:2672:G:C3'	35:DA:2673:G:H5''	2.45	0.46
35:DA:359:A:H2'	35:DA:360:G:C8	2.50	0.46
35:DA:524:U:C2'	35:DA:524:U:O2	2.60	0.46
36:DB:32:C:OP2	41:DG:96:ARG:NH2	2.35	0.46
36:DB:35:U:HO2'	36:DB:36:C:H5'	1.79	0.46
35:DA:2127:G:H5'	37:DC:36:LYS:NZ	2.31	0.46
39:DE:168:MET:O	39:DE:170:LEU:HD12	2.15	0.46
41:DG:142:PRO:O	41:DG:144:ILE:N	2.48	0.46
42:DH:74:ASN:C	42:DH:76:VAL:N	2.69	0.46
44:DN:107:LEU:HB2	44:DN:108:PRO:CD	2.45	0.46
44:DN:41:ASP:O	44:DN:42:TRP:C	2.54	0.46
45:DO:1:MET:HE2	45:DO:1:MET:N	2.29	0.46
35:DA:1202:C:O2	46:DP:7:ARG:NH2	2.48	0.46
47:DQ:103:MET:HE2	47:DQ:125:LEU:HD21	1.98	0.46
47:DQ:70:PRO:CA	47:DQ:95:ALA:HB2	2.46	0.46
48:DR:12:ARG:HG3	48:DR:12:ARG:NH1	2.28	0.46
49:DS:101:LEU:C	49:DS:101:LEU:HD13	2.35	0.46
50:DT:106:SER:C	50:DT:107:ASP:OD1	2.53	0.46
35:DA:996:A:O4'	51:DU:92:ARG:NH2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DV:15:GLU:H	52:DV:18:LEU:HD11	1.81	0.46
52:DV:33:VAL:HA	52:DV:62:LEU:O	2.15	0.46
53:DW:5:ALA:HB3	53:DW:105:VAL:N	2.31	0.46
53:DW:84:ARG:HB2	53:DW:96:ILE:CG2	2.46	0.46
28:D2:23:LYS:CB	54:DX:5:TYR:HE1	2.29	0.46
55:DY:99:CYS:O	55:DY:100:ALA:HB2	2.16	0.46
35:DA:310:A:OP1	55:DY:17:SER:O	2.34	0.46
55:DY:14:LEU:O	55:DY:72:VAL:HG12	2.16	0.46
55:DY:88:LYS:NZ	55:DY:93:GLY:N	2.47	0.46
56:DZ:146:ILE:O	56:DZ:148:ASP:N	2.48	0.46
56:DZ:16:SER:O	56:DZ:20:ARG:HB2	2.15	0.46
1:AA:122:G:H2'	1:AA:123:C:H6	1.80	0.46
1:AA:1231:G:O2'	1:AA:1232:U:H5'	2.15	0.46
1:AA:1271:G:H2'	1:AA:1272:G:O4'	2.15	0.46
1:AA:1511:G:C2'	1:AA:1512:U:H5'	2.45	0.46
1:AA:302:G:H2'	1:AA:303:A:O4'	2.16	0.46
1:AA:670:G:H2'	1:AA:671:G:O4'	2.15	0.46
1:AA:694:A:O2'	23:AW:39:A:C1'	2.61	0.46
1:AA:761:G:H2'	1:AA:762:C:C6	2.51	0.46
1:AA:929:G:C2'	1:AA:930:C:H5'	2.46	0.46
2:AB:114:ARG:HH11	2:AB:118:LEU:HD23	1.81	0.46
2:AB:178:ARG:CB	2:AB:178:ARG:HH11	2.27	0.46
2:AB:167:PRO:CD	2:AB:188:ALA:HB2	2.46	0.46
2:AB:41:ILE:CD1	2:AB:41:ILE:N	2.78	0.46
2:AB:97:TRP:HH2	2:AB:176:GLU:CB	2.27	0.46
3:AC:84:ILE:HG13	3:AC:101:LEU:HD13	1.97	0.46
5:AE:50:GLU:CD	5:AE:51:VAL:H	2.19	0.46
6:AF:66:GLU:O	6:AF:67:MET:HB3	2.15	0.46
7:AG:26:PHE:CZ	7:AG:30:ILE:HD11	2.51	0.46
9:AI:25:LYS:O	9:AI:25:LYS:HG3	2.16	0.46
10:AJ:31:GLY:HA3	10:AJ:78:ASN:ND2	2.31	0.46
13:AM:90:LEU:O	13:AM:91:ARG:HB2	2.16	0.46
14:AN:6:LEU:C	14:AN:8:GLU:N	2.69	0.46
17:AQ:10:VAL:CG1	17:AQ:53:LEU:HA	2.45	0.46
25:AY:70:SER:CB	25:AY:76:LEU:HD12	2.35	0.46
25:AY:92:PRO:HA	25:AY:101:ILE:HG23	1.97	0.46
26:B0:72:ARG:CD	26:B0:75:LEU:HD13	2.42	0.46
29:B3:12:PRO:HA	29:B3:15:TYR:HD1	1.81	0.46
33:B7:15:THR:HG22	33:B7:16:HIS:NE2	2.30	0.46
33:B7:5:TRP:C	33:B7:6:GLN:NE2	2.68	0.46
35:BA:818:G:H3'	35:BA:1187:G:H22	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1305:C:O2'	35:BA:1306:C:H5'	2.15	0.46
35:BA:1435:G:H2'	35:BA:1436:G:O4'	2.16	0.46
35:BA:1490:A:H2	38:BD:75:ILE:HD12	1.81	0.46
35:BA:1550:C:H2'	35:BA:1551:C:C6	2.40	0.46
35:BA:1668:A:C8	35:BA:1674:G:C6	3.04	0.46
35:BA:1820:U:H4'	35:BA:1821:A:OP2	2.15	0.46
35:BA:1839:G:C5	35:BA:1840:G:N7	2.84	0.46
35:BA:201:C:H2'	35:BA:202:U:H5'	1.98	0.46
35:BA:2052:G:C4	35:BA:2053:G:C8	3.03	0.46
35:BA:2079:U:H2'	35:BA:2080:G:H8	1.79	0.46
35:BA:2251:G:H2'	35:BA:2252:G:O4'	2.16	0.46
35:BA:919:G:N2	35:BA:2269:A:OP2	2.44	0.46
35:BA:2276:G:OP2	47:BQ:84:GLY:N	2.49	0.46
35:BA:2556:C:H2'	35:BA:2557:G:O4'	2.16	0.46
35:BA:2640:G:H8	35:BA:2640:G:H5'	1.81	0.46
35:BA:2822:G:H2'	35:BA:2823:A:H5''	1.97	0.46
35:BA:374:A:H2'	35:BA:375:C:O4'	2.15	0.46
35:BA:472:A:H2'	35:BA:473:G:H5'	1.97	0.46
35:BA:641:C:H2'	35:BA:642:G:O4'	2.16	0.46
35:BA:675:A:OP1	40:BF:63:LYS:HD2	2.16	0.46
35:BA:822:U:H2'	35:BA:822:U:O2	2.16	0.46
35:BA:85:G:N2	35:BA:86:C:H1'	2.31	0.46
38:BD:21:PHE:O	38:BD:22:SER:C	2.53	0.46
39:BE:197:ILE:HG13	39:BE:199:ARG:NH1	2.27	0.46
35:BA:442:G:H4'	40:BF:46:ARG:HD3	1.97	0.46
41:BG:114:ILE:CG2	41:BG:115:ARG:N	2.79	0.46
41:BG:140:ILE:HD12	41:BG:141:PHE:N	2.30	0.46
41:BG:14:GLU:O	41:BG:18:GLU:HB3	2.15	0.46
44:BN:62:VAL:O	44:BN:63:THR:CG2	2.62	0.46
45:BO:101:PRO:C	45:BO:102:VAL:HG22	2.35	0.46
45:BO:60:ALA:HB1	45:BO:85:VAL:O	2.15	0.46
46:BP:57:THR:C	46:BP:59:LEU:H	2.17	0.46
46:BP:65:ARG:HG2	46:BP:65:ARG:O	2.16	0.46
46:BP:75:ILE:CD1	46:BP:75:ILE:H	2.08	0.46
47:BQ:108:GLY:HA3	56:BZ:116:VAL:HG11	1.98	0.46
48:BR:13:HIS:C	48:BR:13:HIS:ND1	2.67	0.46
48:BR:17:ARG:O	48:BR:18:LEU:C	2.53	0.46
35:BA:1453:U:OP1	48:BR:77:ARG:NH1	2.49	0.46
50:BT:48:ILE:HG22	50:BT:49:VAL:N	2.31	0.46
51:BU:83:LEU:HG	51:BU:88:ILE:HG12	1.98	0.46
52:BV:39:LEU:HD11	52:BV:53:GLU:H	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:10:VAL:O	53:BW:11:ARG:CB	2.64	0.46
53:BW:44:ALA:O	53:BW:48:ALA:N	2.44	0.46
54:BX:64:LYS:CG	54:BX:65:ARG:H	2.28	0.46
1:CA:116:A:H2'	1:CA:117:G:O4'	2.16	0.46
1:CA:1523:G:C5	1:CA:1524:C:C5	3.04	0.46
1:CA:189(I):G:H2'	1:CA:189(J):G:H8	1.80	0.46
1:CA:23:C:C2'	1:CA:24:U:H5'	2.46	0.46
1:CA:435:C:H2'	1:CA:435:C:O2	2.15	0.46
2:CB:164:VAL:CG1	2:CB:165:VAL:N	2.79	0.46
3:CC:153:VAL:HG12	3:CC:154:SER:H	1.80	0.46
3:CC:159:GLY:O	3:CC:160:ALA:C	2.54	0.46
4:CD:68:TYR:N	4:CD:68:TYR:CD1	2.84	0.46
5:CE:37:ARG:HA	5:CE:112:LEU:O	2.16	0.46
6:CF:14:LEU:HD13	6:CF:15:ASP:O	2.15	0.46
7:CG:39:ALA:O	7:CG:41:ARG:N	2.49	0.46
8:CH:103:VAL:CG1	8:CH:108:GLY:HA3	2.45	0.46
9:CI:89:ASN:C	9:CI:91:ASP:H	2.19	0.46
10:CJ:48:THR:HG23	10:CJ:62:HIS:CA	2.46	0.46
1:CA:303:A:P	12:CL:17:LYS:HE3	2.56	0.46
1:CA:608:A:H4'	16:CP:32:TYR:OH	2.16	0.46
19:CS:15:LEU:H	19:CS:15:LEU:CD2	2.28	0.46
20:CT:64:ASP:C	20:CT:66:ALA:N	2.68	0.46
20:CT:83:ARG:HG2	20:CT:86:ARG:HD3	1.98	0.46
24:CX:17:U:O5'	24:CX:17:U:H6	1.98	0.46
26:D0:29:GLN:HB2	26:D0:67:VAL:HG23	1.97	0.46
27:D1:86:SER:C	27:D1:89:GLU:OE2	2.54	0.46
27:D1:86:SER:CA	27:D1:89:GLU:OE1	2.54	0.46
28:D2:50:ILE:O	28:D2:51:ARG:CB	2.63	0.46
35:DA:83:G:N2	35:DA:103:A:OP2	2.46	0.46
35:DA:1371:G:H8	35:DA:1371:G:O5'	1.98	0.46
35:DA:1428:C:C4	35:DA:1569:A:H5"	2.49	0.46
35:DA:1271:G:N2	35:DA:1617:C:C4'	2.79	0.46
35:DA:1613:G:N1	35:DA:1619:G:C5	2.83	0.46
35:DA:1659:U:OP2	39:DE:132:HIS:HE1	1.99	0.46
35:DA:1714:G:H2'	35:DA:1717:G:C8	2.50	0.46
35:DA:1947:C:O2'	35:DA:1948:G:H5'	2.16	0.46
35:DA:227:A:C2	35:DA:2407:G:C1'	2.97	0.46
35:DA:2701:C:H2'	35:DA:2702:U:H6	1.81	0.46
35:DA:2762:G:C3'	35:DA:2763:G:C5'	2.93	0.46
35:DA:514:A:O2'	35:DA:515:A:H5'	2.14	0.46
35:DA:614(B):G:C4	40:DF:44:ARG:NH2	2.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:733:G:O6	35:DA:761:A:N7	2.49	0.46
35:DA:80:G:O2'	35:DA:81:G:H5'	2.15	0.46
35:DA:954:G:OP1	47:DQ:15:GLY:N	2.48	0.46
39:DE:1:MET:HE2	39:DE:1:MET:H1	1.81	0.46
40:DF:101:LEU:HD12	40:DF:102:PRO:CD	2.36	0.46
40:DF:170:LEU:HD21	40:DF:172:TRP:CZ2	2.51	0.46
41:DG:128:ARG:O	41:DG:129:GLY:O	2.33	0.46
41:DG:36:LYS:HB3	41:DG:160:VAL:CG2	2.45	0.46
41:DG:41:GLN:CG	41:DG:43:LEU:HD12	2.27	0.46
43:DI:114:LEU:O	43:DI:115:ALA:CB	2.64	0.46
44:DN:12:ARG:NH2	44:DN:39:ARG:NH1	2.63	0.46
44:DN:22:THR:HA	44:DN:61:ARG:CB	2.33	0.46
44:DN:34:LEU:HA	44:DN:34:LEU:HD22	1.77	0.46
44:DN:32:THR:HG23	44:DN:37:LYS:HB3	1.96	0.46
44:DN:94:HIS:N	44:DN:94:HIS:ND1	2.63	0.46
46:DP:14:LYS:O	46:DP:15:ARG:CB	2.64	0.46
46:DP:59:LEU:HA	46:DP:61:ARG:CD	2.45	0.46
49:DS:83:LYS:O	49:DS:85:VAL:N	2.48	0.46
49:DS:92:TYR:HD1	49:DS:92:TYR:C	2.18	0.46
52:DV:5:VAL:HB	52:DV:37:VAL:O	2.15	0.46
52:DV:83:ARG:HH11	52:DV:83:ARG:CG	2.08	0.46
53:DW:48:ALA:O	53:DW:49:LYS:C	2.53	0.46
55:DY:96:ILE:HG13	55:DY:100:ALA:H	1.80	0.46
56:DZ:8:TYR:CD1	56:DZ:8:TYR:N	2.83	0.46
1:AA:1001:A:H2	1:AA:1001(A):G:O6	1.98	0.46
1:AA:1011:G:O2'	1:AA:1012:U:H5'	2.16	0.46
1:AA:1049:U:O2'	1:AA:1050:G:OP2	2.29	0.46
1:AA:1097:C:H2'	1:AA:1098:C:H6	1.79	0.46
1:AA:1405:G:H1	1:AA:1496:C:H42	1.64	0.46
1:AA:197:A:N6	1:AA:221:C:H5'	2.31	0.46
1:AA:374:A:C6	1:AA:375:U:C4	3.04	0.46
1:AA:726:C:H2'	1:AA:727:G:H8	1.81	0.46
1:AA:895:G:C6	1:AA:896:C:N4	2.84	0.46
2:AB:173:ALA:O	2:AB:174:VAL:C	2.54	0.46
3:AC:150:LYS:O	3:AC:201:TYR:N	2.45	0.46
4:AD:13:ARG:NH1	4:AD:40:PRO:HA	2.30	0.46
5:AE:20:GLN:O	5:AE:23:GLY:O	2.34	0.46
6:AF:54:LYS:O	6:AF:56:PRO:HD3	2.15	0.46
7:AG:42:ILE:HG23	7:AG:117:ALA:HA	1.97	0.46
8:AH:114:THR:CG2	8:AH:117:GLY:O	2.64	0.46
8:AH:127:LEU:HD12	8:AH:129:VAL:HG13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:49:GLU:HG2	8:AH:62:TYR:HE2	1.81	0.46
8:AH:51:VAL:HG11	8:AH:60:ARG:CG	2.46	0.46
9:AI:89:ASN:HB3	9:AI:92:TYR:CD1	2.51	0.46
10:AJ:3:LYS:O	10:AJ:100:THR:HA	2.15	0.46
11:AK:102:GLY:O	11:AK:103:LEU:O	2.34	0.46
11:AK:86:GLY:N	11:AK:112:THR:OG1	2.48	0.46
13:AM:97:PRO:O	13:AM:98:VAL:N	2.49	0.46
18:AR:72:ARG:O	18:AR:75:ILE:N	2.43	0.46
18:AR:87:ARG:HH11	18:AR:87:ARG:HB3	1.77	0.46
19:AS:40:ILE:HD13	19:AS:62:ILE:CD1	2.45	0.46
20:AT:74:LYS:CG	20:AT:75:ASN:H	2.28	0.46
22:AV:37:A:C8	22:AV:38:U:C5	3.04	0.46
23:AW:38:A:H2'	23:AW:39:A:H5'	1.98	0.46
25:AY:29:ARG:NH1	25:AY:110:ARG:NH2	2.64	0.46
25:AY:149:LEU:HB3	25:AY:153:GLU:CB	2.46	0.46
27:B1:37:ILE:CD1	27:B1:37:ILE:H	2.11	0.46
31:B5:42:PRO:C	31:B5:43:HIS:HD2	2.20	0.46
35:BA:109:G:O2'	35:BA:110:G:H5'	2.15	0.46
35:BA:1223:G:C5	35:BA:1225:G:OP2	2.68	0.46
35:BA:1259:G:H2'	35:BA:1260:G:C8	2.50	0.46
35:BA:1426:G:C6	35:BA:1427:A:N1	2.84	0.46
35:BA:1553:A:N7	35:BA:1555:G:C5	2.83	0.46
35:BA:1662:C:O2'	35:BA:2687:U:OP1	2.34	0.46
35:BA:1665:A:C2'	35:BA:1666:G:H5'	2.46	0.46
35:BA:1800:C:C4	35:BA:1818:U:O2	2.69	0.46
35:BA:2008:C:H2'	35:BA:2009:G:C8	2.51	0.46
35:BA:2193:G:C4	35:BA:2194:G:C8	3.04	0.46
35:BA:2194:G:C4	35:BA:2195:C:C5	3.04	0.46
35:BA:2206:G:N2	35:BA:2207:G:C5'	2.73	0.46
35:BA:2220:G:C4	35:BA:2221:G:C8	3.03	0.46
35:BA:2355:C:C4	35:BA:2356:C:C4	3.04	0.46
35:BA:2473:U:O2	35:BA:2473:U:C2'	2.64	0.46
35:BA:2571:C:H5'	35:BA:2572:A:C5'	2.46	0.46
35:BA:2577:A:H5''	35:BA:2578:G:C5'	2.42	0.46
35:BA:2714:G:C5	35:BA:2715:C:C4	3.04	0.46
35:BA:2842:G:C6	35:BA:2876:G:C6	3.04	0.46
35:BA:286:C:C2'	35:BA:286:C:O2	2.60	0.46
35:BA:2883:A:H5'	35:BA:2884:U:H5'	1.98	0.46
35:BA:304:G:H1	35:BA:313:C:N4	2.11	0.46
35:BA:433:C:H2'	35:BA:434:U:C6	2.51	0.46
35:BA:447:A:N3	35:BA:473:G:C8	2.84	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:701:G:O2'	35:BA:702:G:H5'	2.15	0.46
35:BA:737:C:C2'	35:BA:738:G:O5'	2.63	0.46
35:BA:811:U:H1'	35:BA:1251:C:H5''	1.96	0.46
37:BC:51:PRO:O	37:BC:52:ARG:HB2	2.16	0.46
39:BE:120:TRP:O	39:BE:121:ASN:C	2.53	0.46
39:BE:9:VAL:HG23	50:BT:8:LYS:HB2	1.97	0.46
41:BG:113:ARG:O	41:BG:114:ILE:CG1	2.64	0.46
41:BG:169:ALA:C	41:BG:173:LEU:HD23	2.37	0.46
41:BG:27:ASN:ND2	41:BG:29:TRP:H	2.14	0.46
42:BH:25:LYS:HB3	42:BH:32:GLU:OE2	2.15	0.46
43:BI:94:ALA:HA	43:BI:97:ILE:HG13	1.96	0.46
44:BN:13:TRP:O	44:BN:135:PRO:HD2	2.15	0.46
45:BO:43:VAL:HG11	45:BO:46:ALA:HB2	1.98	0.46
46:BP:16:ARG:O	46:BP:18:ARG:N	2.49	0.46
49:BS:56:LEU:O	49:BS:57:LYS:HB3	2.16	0.46
50:BT:31:SER:O	50:BT:32:TYR:O	2.34	0.46
35:BA:1011:G:OP2	51:BU:70:ARG:NH2	2.49	0.46
53:BW:73:ALA:HB3	53:BW:106:ILE:CD1	2.28	0.46
54:BX:83:VAL:O	54:BX:83:VAL:HG23	2.16	0.46
56:BZ:128:VAL:HG22	56:BZ:129:SER:N	2.31	0.46
56:BZ:5:LEU:HB2	56:BZ:59:LEU:CD2	2.45	0.46
1:CA:1002:G:H2'	1:CA:1003:G:O4'	2.16	0.46
1:CA:1095:U:H5'	1:CA:1109:C:O2	2.15	0.46
1:CA:1111:A:N1	3:CC:177:THR:OG1	2.44	0.46
1:CA:1227:A:O2'	13:CM:115:LYS:HB2	2.16	0.46
1:CA:1236:A:H4'	1:CA:1304:G:H4'	1.97	0.46
1:CA:1423:G:N3	1:CA:1424:C:C6	2.84	0.46
1:CA:377:G:O2'	1:CA:378:G:H5'	2.15	0.46
1:CA:411:A:H2'	1:CA:412:A:H4'	1.97	0.46
1:CA:544:G:H2'	1:CA:545:C:H6	1.79	0.46
1:CA:577:G:C8	1:CA:816:A:C6	3.04	0.46
1:CA:62:U:H5''	1:CA:385:C:O2	2.16	0.46
1:CA:651:C:H2'	1:CA:652:U:H6	1.81	0.46
1:CA:678:U:H3	1:CA:713:G:N2	2.14	0.46
1:CA:778:G:H2'	1:CA:779:C:O4'	2.15	0.46
1:CA:789:U:H2'	1:CA:789:U:O2	2.16	0.46
1:CA:794:A:O2'	1:CA:795:C:H5'	2.16	0.46
3:CC:87:LEU:CB	3:CC:101:LEU:HD11	2.46	0.46
3:CC:109:PRO:HA	3:CC:115:LEU:HD12	1.97	0.46
4:CD:148:VAL:HG21	4:CD:181:MET:HB2	1.98	0.46
4:CD:30:LYS:HA	4:CD:35:ARG:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1118:C:P	9:CI:104:ARG:HG3	2.56	0.46
1:CA:1295:G:O2'	13:CM:14:ARG:NH1	2.48	0.46
13:CM:13:LYS:O	13:CM:45:VAL:HG23	2.16	0.46
15:CO:66:LEU:O	15:CO:67:LEU:C	2.54	0.46
17:CQ:12:SER:HA	17:CQ:14:LYS:NZ	2.30	0.46
17:CQ:29:HIS:CE1	17:CQ:31:LEU:H	2.33	0.46
17:CQ:82:MET:HA	17:CQ:85:VAL:HG23	1.97	0.46
25:CY:79:ILE:O	25:CY:82:ALA:HB3	2.16	0.46
26:D0:36:ILE:HA	26:D0:60:PHE:HB3	1.97	0.46
35:DA:1406:U:C3'	35:DA:1407:C:H6	2.29	0.46
35:DA:1946:U:HO2'	35:DA:1947:C:H5'	1.81	0.46
35:DA:210:C:H2'	35:DA:211:A:C8	2.51	0.46
35:DA:2120:G:O2'	35:DA:2121:G:H5'	2.16	0.46
35:DA:2261:C:C2	35:DA:2280:G:C2	3.04	0.46
35:DA:2419:U:H2'	35:DA:2420:C:H6	1.80	0.46
35:DA:2773:C:OP1	39:DE:164:ARG:HG2	2.15	0.46
35:DA:2792:G:N3	35:DA:2792:G:H2'	2.30	0.46
35:DA:2865:U:H3'	35:DA:2866:U:O2	2.16	0.46
35:DA:451:C:H41	35:DA:453:C:H3'	1.80	0.46
35:DA:661:C:H2'	35:DA:662:G:H8	1.77	0.46
35:DA:764:A:N1	35:DA:781:A:C2	2.84	0.46
35:DA:884:C:H4'	35:DA:892:G:N7	2.30	0.46
38:DD:123:ALA:HB3	38:DD:131:LEU:HD23	1.98	0.46
38:DD:16:MET:HG3	38:DD:206:LEU:O	2.15	0.46
40:DF:51:THR:HG23	40:DF:92:PRO:HG2	1.98	0.46
41:DG:114:ILE:HG13	41:DG:117:PHE:CB	2.46	0.46
41:DG:144:ILE:HD12	41:DG:145:THR:N	2.28	0.46
42:DH:130:ARG:CB	42:DH:130:ARG:HH11	2.29	0.46
44:DN:29:LYS:C	44:DN:31:ALA:H	2.19	0.46
45:DO:64:ARG:HD3	45:DO:101:PRO:C	2.36	0.46
45:DO:62:VAL:CG1	45:DO:65:THR:HG22	2.46	0.46
45:DO:59:LYS:O	45:DO:86:ILE:HG23	2.16	0.46
35:DA:2415:G:C4'	46:DP:66:GLY:HA3	2.44	0.46
47:DQ:53:ALA:CA	47:DQ:56:ARG:HB3	2.45	0.46
35:DA:2839:G:H1'	48:DR:93:GLY:H	1.80	0.46
49:DS:101:LEU:HD21	49:DS:103:GLU:HG3	1.97	0.46
49:DS:95:HIS:O	49:DS:98:VAL:HG23	2.16	0.46
52:DV:82:ARG:NH1	52:DV:84:LYS:HD3	2.31	0.46
53:DW:66:GLU:HA	53:DW:69:LEU:HD21	1.96	0.46
54:DX:37:THR:HG23	54:DX:54:VAL:HB	1.98	0.46
1:AA:1396:A:H4'	1:AA:1398:A:H1'	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:29:G:H5'	1:AA:296:U:OP1	2.16	0.46
1:AA:116:A:H61	1:AA:313:A:H1'	1.80	0.46
1:AA:517:G:N3	1:AA:531:U:H5'	2.31	0.46
1:AA:774:G:H1	1:AA:805:C:N4	2.13	0.46
1:AA:818:G:C3'	1:AA:819:A:C5'	2.91	0.46
1:AA:819:A:C8	1:AA:1529:G:N1	2.83	0.46
1:AA:854:G:OP2	1:AA:871:U:C5	2.69	0.46
1:AA:1075:C:H5'	2:AB:103:THR:HG21	1.98	0.46
2:AB:165:VAL:CG2	2:AB:166:ASP:N	2.58	0.46
2:AB:36:ARG:HB3	2:AB:41:ILE:CD1	2.42	0.46
3:AC:178:LEU:H	3:AC:178:LEU:HD22	1.81	0.46
4:AD:13:ARG:CG	4:AD:14:ARG:H	2.28	0.46
4:AD:13:ARG:HG2	4:AD:14:ARG:H	1.81	0.46
4:AD:24:GLU:O	4:AD:27:TYR:HB2	2.15	0.46
4:AD:65:ARG:NH1	4:AD:72:GLU:CA	2.78	0.46
1:AA:1080:A:C5'	5:AE:16:THR:HG21	2.46	0.46
1:AA:932:C:H4'	7:AG:4:ARG:NH2	2.31	0.46
11:AK:73:MET:SD	11:AK:103:LEU:HD13	2.56	0.46
7:AG:149:ARG:HB3	11:AK:59:TYR:CE2	2.51	0.46
11:AK:60:ALA:O	11:AK:61:ALA:C	2.54	0.46
11:AK:72:ALA:HB1	11:AK:77:MET:HE3	1.97	0.46
11:AK:73:MET:HA	11:AK:77:MET:HB3	1.98	0.46
12:AL:104:VAL:HG12	12:AL:105:TYR:N	2.30	0.46
13:AM:91:ARG:HB2	13:AM:98:VAL:HG22	1.96	0.46
14:AN:3:ARG:O	14:AN:7:ILE:HG23	2.16	0.46
19:AS:36:ARG:HB3	19:AS:36:ARG:NH1	2.31	0.46
20:AT:98:PRO:O	20:AT:100:ILE:N	2.49	0.46
25:AY:114:LEU:CD2	25:AY:183:ILE:HG23	2.46	0.46
25:AY:32:ARG:HH21	25:AY:88:LEU:HG	1.81	0.46
25:AY:70:SER:HB3	25:AY:76:LEU:HB2	1.97	0.46
27:B1:47:GLN:NE2	27:B1:47:GLN:C	2.69	0.46
29:B3:11:SER:OG	29:B3:12:PRO:HD2	2.16	0.46
33:B7:24:THR:O	33:B7:28:ARG:HG3	2.16	0.46
34:B8:52:LYS:HE3	34:B8:52:LYS:CA	2.43	0.46
35:BA:1021:A:C3'	35:BA:1021:A:C8	2.99	0.46
35:BA:1130:U:O2	35:BA:2025:C:H5''	2.15	0.46
35:BA:1152:C:O2'	35:BA:1153:C:H5'	2.15	0.46
35:BA:552:G:H1'	35:BA:1220:A:C2	2.51	0.46
35:BA:1281:G:H2'	35:BA:1282:U:H6	1.81	0.46
35:BA:1331:A:H2'	35:BA:1333:C:H5	1.81	0.46
35:BA:134:C:O2'	35:BA:135:G:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1467:C:N4	35:BA:1525:G:H1	2.14	0.46
35:BA:1387:C:C5'	35:BA:1469:A:H4'	2.46	0.46
35:BA:1479:G:H2'	35:BA:1480:G:O4'	2.15	0.46
35:BA:1819:A:H5''	38:BD:161:THR:CG2	2.46	0.46
35:BA:1831:G:H2'	35:BA:1832:C:H6	1.81	0.46
35:BA:1853:A:N1	35:BA:2087:G:H1'	2.30	0.46
35:BA:2362:G:O2'	35:BA:2363:C:H5'	2.15	0.46
35:BA:2598:A:H5''	38:BD:236:GLY:CA	2.42	0.46
35:BA:2605:U:H2'	35:BA:2606:C:C6	2.51	0.46
35:BA:746:A:C5	35:BA:2611:U:H5''	2.51	0.46
35:BA:2845:G:H5''	50:BT:55:ASN:HA	1.98	0.46
35:BA:467:G:H2'	35:BA:468:G:C8	2.40	0.46
36:BB:66:A:C2	36:BB:109:C:C2	3.04	0.46
38:BD:117:VAL:CG2	38:BD:118:VAL:N	2.75	0.46
38:BD:172:TYR:CD1	38:BD:186:HIS:CA	2.94	0.46
39:BE:179:GLU:O	39:BE:180:ASN:HB3	2.16	0.46
39:BE:83:ASP:O	39:BE:84:PHE:HB2	2.16	0.46
40:BF:167:ALA:HB1	40:BF:173:VAL:CG1	2.46	0.46
41:BG:68:PRO:HA	41:BG:92:VAL:HB	1.98	0.46
42:BH:85:LYS:HD3	42:BH:133:VAL:HB	1.98	0.46
42:BH:43:VAL:HB	42:BH:51:ARG:O	2.16	0.46
42:BH:19:VAL:CG1	42:BH:44:VAL:HG22	2.46	0.46
43:BI:42:SER:C	43:BI:44:LEU:N	2.69	0.46
43:BI:85:GLU:O	43:BI:86:THR:O	2.34	0.46
45:BO:19:ILE:HG22	45:BO:42:SER:C	2.36	0.46
45:BO:14:THR:HG21	45:BO:86:ILE:CG1	2.46	0.46
46:BP:101:VAL:O	46:BP:103:ALA:N	2.49	0.46
46:BP:131:SER:O	46:BP:135:LEU:N	2.49	0.46
47:BQ:20:ALA:HA	47:BQ:98:LYS:HD3	1.98	0.46
48:BR:17:ARG:CG	48:BR:17:ARG:NH1	2.76	0.46
48:BR:38:VAL:CB	48:BR:39:PRO:HD3	2.34	0.46
49:BS:25:ARG:HD3	49:BS:42:ASP:OD2	2.16	0.46
49:BS:87:PHE:CZ	49:BS:97:ARG:NH2	2.83	0.46
50:BT:17:THR:O	50:BT:18:ASP:HB3	2.16	0.46
50:BT:52:ILE:CG2	50:BT:61:PHE:HB2	2.44	0.46
53:BW:14:PRO:O	53:BW:15:ARG:C	2.54	0.46
53:BW:28:SER:O	53:BW:29:LEU:C	2.52	0.46
55:BY:42:VAL:CB	55:BY:65:ALA:HB3	2.46	0.46
55:BY:84:ARG:O	55:BY:85:VAL:HG22	2.16	0.46
56:BZ:127:LYS:HB2	56:BZ:164:ALA:HB2	1.96	0.46
56:BZ:47:VAL:O	56:BZ:51:ALA:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1349:A:H3'	9:CI:118:LYS:CE	2.34	0.46
1:CA:16:A:C2	1:CA:17:U:C5	3.04	0.46
1:CA:182:U:H2'	1:CA:183:G:O4'	2.16	0.46
1:CA:136:C:H42	1:CA:227:G:H1	1.64	0.46
1:CA:343:U:O2'	1:CA:344:A:H2'	2.16	0.46
1:CA:474:G:O2'	1:CA:475:G:H5'	2.16	0.46
1:CA:685:G:H21	1:CA:686:U:H3	1.63	0.46
1:CA:728:A:C2	1:CA:729:A:C5	3.04	0.46
1:CA:829:G:H2'	1:CA:830:G:C8	2.51	0.46
1:CA:893:C:H2'	1:CA:894:G:H8	1.81	0.46
1:CA:918:A:H2'	1:CA:919:A:C8	2.51	0.46
2:CB:178:ARG:HH11	2:CB:178:ARG:CG	2.28	0.46
2:CB:54:THR:HG21	2:CB:201:ILE:CD1	2.44	0.46
2:CB:83:MET:O	2:CB:86:GLU:N	2.49	0.46
1:CA:407:G:O2'	4:CD:116:GLN:HG3	2.16	0.46
4:CD:61:LYS:HA	4:CD:203:VAL:CG2	2.31	0.46
5:CE:101:ILE:HG12	5:CE:118:ILE:O	2.16	0.46
5:CE:147:ASP:O	5:CE:151:LEU:HG	2.16	0.46
9:CI:28:VAL:HG22	9:CI:63:ILE:O	2.15	0.46
9:CI:4:TYR:HA	9:CI:88:TYR:CE1	2.51	0.46
16:CP:23:ASP:OD2	16:CP:25:ARG:NH2	2.49	0.46
20:CT:26:ASN:O	20:CT:27:LYS:C	2.54	0.46
23:CW:34:U:O2'	23:CW:36:A:N7	2.45	0.46
25:CY:3:LEU:O	25:CY:4:LYS:C	2.53	0.46
28:D2:15:LYS:O	28:D2:17:SER:N	2.49	0.46
31:D5:42:PRO:C	31:D5:43:HIS:CD2	2.89	0.46
34:D8:52:LYS:HE3	34:D8:52:LYS:CA	2.43	0.46
35:DA:105:C:H5'	35:DA:106:C:OP2	2.15	0.46
35:DA:1214:A:H2'	35:DA:1215:G:C8	2.51	0.46
35:DA:1360:A:C5	35:DA:1372:U:C4	3.04	0.46
35:DA:1644:C:C2'	35:DA:1645:G:H5'	2.46	0.46
35:DA:1710:C:O2'	35:DA:1711:C:H5'	2.16	0.46
35:DA:1916:A:H5'	35:DA:1917:U:OP2	2.15	0.46
35:DA:571:A:H5''	35:DA:2030:A:H62	1.78	0.46
35:DA:2193:G:C4	35:DA:2194:G:C8	3.04	0.46
35:DA:2240:C:O2'	35:DA:2241:A:H5'	2.16	0.46
35:DA:2263:C:O2'	35:DA:2264:C:H5'	2.16	0.46
35:DA:2312:U:H2'	35:DA:2313:C:C5'	2.43	0.46
35:DA:2732:G:H2'	35:DA:2733:A:H5'	1.96	0.46
35:DA:374:A:H2'	35:DA:375:C:H5'	1.98	0.46
35:DA:394:A:H2'	35:DA:395:U:H5'	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:790:C:H4'	35:DA:790:C:OP1	2.16	0.46
35:DA:995:C:C2	51:DU:57:PHE:HE2	2.33	0.46
36:DB:63:G:N3	36:DB:63:G:H2'	2.31	0.46
38:DD:211:ARG:NH1	38:DD:211:ARG:HG2	2.31	0.46
39:DE:115:GLY:HA2	39:DE:157:ALA:CB	2.45	0.46
40:DF:9:ILE:HG23	40:DF:14:PRO:HA	1.98	0.46
41:DG:131:TYR:O	41:DG:159:VAL:CG2	2.64	0.46
41:DG:14:GLU:HG2	41:DG:15:VAL:N	2.30	0.46
41:DG:35:GLU:HB3	41:DG:160:VAL:HG11	1.98	0.46
43:DI:69:LYS:C	43:DI:71:ILE:H	2.19	0.46
44:DN:46:VAL:CG1	44:DN:47:ALA:H	1.99	0.46
47:DQ:28:ALA:HB3	47:DQ:105:GLU:OE2	2.15	0.46
47:DQ:50:ALA:O	47:DQ:54:MET:CB	2.64	0.46
48:DR:88:ARG:NH2	48:DR:89:ASP:OD1	2.48	0.46
50:DT:113:LYS:C	50:DT:114:LEU:HD23	2.35	0.46
50:DT:53:ARG:NE	50:DT:60:THR:OG1	2.49	0.46
55:DY:46:LYS:C	55:DY:47:LYS:HD2	2.36	0.46
56:DZ:110:GLY:O	56:DZ:111:VAL:HG12	2.16	0.46
56:DZ:53:ILE:HG22	56:DZ:71:VAL:HG23	1.97	0.46
1:AA:1016:A:H2'	1:AA:1017:G:H5'	1.97	0.46
1:AA:1030(C):G:N7	1:AA:1031:G:N2	2.64	0.46
1:AA:106:C:O2'	1:AA:107:G:H5'	2.16	0.46
1:AA:1333:A:H2'	1:AA:1334:G:O4'	2.16	0.46
1:AA:1466:C:H2'	1:AA:1467:G:C8	2.51	0.46
1:AA:544:G:C4	1:AA:545:C:C5	3.03	0.46
1:AA:777:A:H2'	1:AA:778:G:C8	2.51	0.46
1:AA:984:C:H2'	1:AA:985:C:C6	2.51	0.46
2:AB:115:LEU:HD21	2:AB:153:ARG:NE	2.31	0.46
2:AB:69:LEU:HB2	2:AB:162:ILE:HG22	1.98	0.46
3:AC:84:ILE:CA	3:AC:87:LEU:HD12	2.40	0.46
7:AG:64:GLN:O	7:AG:66:VAL:N	2.49	0.46
7:AG:84:ASN:HB2	23:AW:33:C:O2'	2.16	0.46
9:AI:99:LEU:O	9:AI:100:GLY:C	2.54	0.46
1:AA:972:C:C4'	10:AJ:57:LYS:HG3	2.44	0.46
12:AL:90:VAL:CG1	12:AL:93:LEU:HG	2.44	0.46
13:AM:89:GLY:O	13:AM:90:LEU:O	2.34	0.46
16:AP:19:ILE:HB	16:AP:37:GLY:C	2.35	0.46
20:AT:43:LEU:HA	20:AT:46:GLU:HB3	1.97	0.46
23:AW:16:C:H4'	23:AW:20:G:OP1	2.16	0.46
23:AW:40:C:H2'	23:AW:41:C:H6	1.80	0.46
25:AY:61:PRO:HD2	25:AY:65:THR:C	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:AY:6:LEU:O	25:AY:9:GLU:N	2.48	0.46
25:AY:84:ARG:C	25:AY:86:SER:N	2.69	0.46
27:B1:46:LEU:N	27:B1:46:LEU:CD1	2.74	0.46
27:B1:48:LYS:HA	27:B1:48:LYS:HE3	1.97	0.46
28:B2:51:ARG:HE	28:B2:55:ARG:NH1	2.14	0.46
29:B3:29:ARG:O	29:B3:30:ARG:C	2.54	0.46
33:B7:16:HIS:HD1	33:B7:21:ARG:NH2	2.14	0.46
35:BA:1131:G:C2	35:BA:1132:A:C5	3.04	0.46
35:BA:118:A:H1'	35:BA:178:G:O4'	2.16	0.46
35:BA:1288:U:C2	35:BA:1327:C:O2	2.69	0.46
35:BA:1444:G:N2	35:BA:1548:C:C2	2.83	0.46
35:BA:1578:U:C2'	35:BA:1579:A:H5''	2.46	0.46
35:BA:1271:G:N2	35:BA:1617:C:C4'	2.79	0.46
35:BA:1653:G:O2'	35:BA:1654:A:OP2	2.28	0.46
35:BA:1675:C:C2	39:BE:129:HIS:CD2	3.04	0.46
35:BA:1803:A:C8	35:BA:1804:C:C5	3.03	0.46
1:AA:1494:G:C5'	35:BA:1913:A:C6	2.99	0.46
35:BA:2052:G:N2	39:BE:149:ARG:HA	2.30	0.46
35:BA:2073:C:H2'	35:BA:2074:U:H6	1.81	0.46
35:BA:2287:A:C2	35:BA:2289:G:C8	3.04	0.46
35:BA:2637:U:H1'	35:BA:2782:G:N2	2.30	0.46
35:BA:2684:U:C2'	35:BA:2685:G:H5'	2.46	0.46
35:BA:2884:U:C2'	35:BA:2885:C:H5'	2.46	0.46
35:BA:372:G:O2'	35:BA:373:U:P	2.74	0.46
35:BA:58:G:H1	35:BA:69:C:N4	2.14	0.46
35:BA:627:A:H8	35:BA:627:A:OP1	1.99	0.46
35:BA:598:G:C6	35:BA:660:G:C6	3.04	0.46
35:BA:774:A:HO2'	35:BA:775:G:H8	1.63	0.46
35:BA:926:A:C4	35:BA:927:G:C8	3.04	0.46
35:BA:979:G:N2	35:BA:985:C:N4	2.64	0.46
36:BB:78:A:H2'	36:BB:79:C:O4'	2.16	0.46
38:BD:233:HIS:O	38:BD:234:GLY:C	2.54	0.46
35:BA:2820:A:C8	39:BE:191:PRO:CB	2.99	0.46
39:BE:4:ILE:HG12	39:BE:28:ALA:HB1	1.96	0.46
39:BE:93:VAL:C	39:BE:95:ILE:N	2.67	0.46
40:BF:132:VAL:O	40:BF:133:ASN:C	2.54	0.46
40:BF:53:THR:HG23	40:BF:56:GLU:CB	2.28	0.46
35:BA:588:U:H1'	40:BF:90:PHE:HB3	1.97	0.46
41:BG:37:VAL:HG12	41:BG:94:LEU:HD12	1.98	0.46
42:BH:92:ILE:O	42:BH:94:TYR:N	2.48	0.46
43:BI:110:ASP:O	43:BI:114:LEU:HG	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BN:122:VAL:HG12	44:BN:123:TYR:N	2.30	0.46
44:BN:28:THR:HG22	44:BN:29:LYS:N	2.31	0.46
44:BN:17:ASP:OD1	44:BN:56:ASN:HB3	2.16	0.46
45:BO:31:LYS:HD2	45:BO:32:TYR:CE1	2.49	0.46
35:BA:1666:G:O3'	45:BO:6:THR:HG23	2.14	0.46
46:BP:75:ILE:HD13	46:BP:77:ARG:NH2	2.30	0.46
48:BR:103:ARG:HB3	48:BR:109:ALA:O	2.15	0.46
35:BA:814:C:OP1	52:BV:84:LYS:HA	2.16	0.46
53:BW:4:LYS:HA	53:BW:106:ILE:HA	1.98	0.46
53:BW:17:VAL:O	53:BW:20:VAL:CG2	2.64	0.46
54:BX:18:TYR:O	54:BX:21:PHE:HB2	2.15	0.46
28:B2:26:ARG:NH2	54:BX:6:ASP:HA	2.31	0.46
55:BY:76:CYS:HB3	55:BY:96:ILE:CD1	2.44	0.46
56:BZ:140:ASP:OD2	56:BZ:140:ASP:N	2.49	0.46
56:BZ:94:GLU:HA	56:BZ:95:PRO:HD2	1.83	0.46
1:CA:1019:C:H2'	1:CA:1020:U:O4'	2.16	0.46
1:CA:1243:C:OP2	21:CU:10:ARG:NH1	2.49	0.46
1:CA:1305:G:H5'	21:CU:4:GLY:CA	2.39	0.46
1:CA:374:A:C6	1:CA:375:U:C4	3.04	0.46
1:CA:434:U:H2'	1:CA:435:C:N1	2.31	0.46
1:CA:583:A:H2'	1:CA:584:G:C8	2.51	0.46
1:CA:670:G:H2'	1:CA:671:G:O4'	2.16	0.46
1:CA:769:G:H1	1:CA:810:C:H42	1.64	0.46
2:CB:142:LEU:HD23	2:CB:142:LEU:C	2.36	0.46
2:CB:142:LEU:O	2:CB:142:LEU:HD23	2.16	0.46
2:CB:221:LEU:HD13	2:CB:221:LEU:C	2.37	0.46
3:CC:172:ARG:NH1	3:CC:174:PRO:HG2	2.31	0.46
3:CC:88:ARG:HG2	3:CC:101:LEU:CB	2.45	0.46
3:CC:94:LEU:HD12	3:CC:95:THR:N	2.31	0.46
4:CD:167:GLY:O	4:CD:168:ARG:C	2.55	0.46
4:CD:72:GLU:O	4:CD:76:ARG:N	2.44	0.46
5:CE:147:ASP:C	5:CE:150:ARG:HB3	2.37	0.46
8:CH:26:VAL:CG2	8:CH:32:LYS:NZ	2.73	0.46
8:CH:53:VAL:HG12	8:CH:54:ASP:OD2	2.16	0.46
9:CI:27:THR:C	9:CI:28:VAL:HG23	2.37	0.46
9:CI:37:PHE:CZ	9:CI:74:ILE:HG12	2.51	0.46
9:CI:99:LEU:O	9:CI:100:GLY:C	2.54	0.46
12:CL:27:LEU:C	12:CL:29:GLY:N	2.69	0.46
15:CO:39:LEU:O	15:CO:42:HIS:HB3	2.15	0.46
17:CQ:15:MET:HG2	17:CQ:16:GLN:N	2.31	0.46
18:CR:37:VAL:O	18:CR:40:LEU:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CT:98:PRO:O	20:CT:100:ILE:N	2.49	0.46
23:CW:14:A:H2'	23:CW:15:G:H5'	1.98	0.46
25:CY:107:THR:HA	25:CY:111:ARG:NH1	2.31	0.46
25:CY:107:THR:OG1	25:CY:108:GLU:N	2.48	0.46
29:D3:7:LYS:O	29:D3:9:VAL:HG13	2.16	0.46
33:D7:48:LYS:CD	33:D7:48:LYS:N	2.78	0.46
35:DA:1164:G:C6	35:DA:1165:U:C4	3.03	0.46
35:DA:1494:A:C3'	35:DA:1494:A:N3	2.73	0.46
35:DA:1494:A:O2'	35:DA:1495:A:H5''	2.16	0.46
35:DA:1346:G:H1	35:DA:1600:C:H42	1.65	0.46
35:DA:1692:U:H2'	35:DA:1694:C:C5	2.51	0.46
35:DA:198:C:H6	35:DA:198:C:O5'	1.99	0.46
35:DA:2008:C:H2'	35:DA:2009:G:C8	2.50	0.46
35:DA:2087:G:C2'	35:DA:2088:G:H5'	2.46	0.46
35:DA:2205:C:H1'	35:DA:2220:G:N2	2.31	0.46
35:DA:2068:U:N3	35:DA:2430:A:C2	2.80	0.46
35:DA:2533:A:C3'	35:DA:2534:A:H5''	2.44	0.46
35:DA:2633:G:N2	35:DA:2634:G:H1'	2.31	0.46
35:DA:2683:C:O2'	35:DA:2684:U:H5'	2.16	0.46
35:DA:49:A:OP1	35:DA:50:U:H3'	2.16	0.46
35:DA:954:G:N3	35:DA:954:G:H2'	2.30	0.46
37:DC:35:ALA:O	37:DC:36:LYS:HE3	2.15	0.46
38:DD:61:LEU:HD12	38:DD:62:TYR:N	2.31	0.46
40:DF:112:MET:O	40:DF:113:ALA:C	2.55	0.46
40:DF:126:VAL:O	40:DF:127:GLU:HB2	2.15	0.46
42:DH:162:ILE:C	42:DH:163:TYR:CD1	2.89	0.46
43:DI:119:PRO:O	43:DI:121:LYS:N	2.48	0.46
43:DI:29:TYR:CD2	43:DI:30:LEU:HD23	2.51	0.46
43:DI:28:ASN:C	43:DI:32:PRO:HG2	2.35	0.46
44:DN:17:ASP:OD1	44:DN:56:ASN:HB3	2.16	0.46
45:DO:68:GLU:HB3	45:DO:78:ARG:CD	2.46	0.46
46:DP:126:VAL:CA	46:DP:145:PRO:HG2	2.45	0.46
46:DP:64:LYS:HB3	46:DP:65:ARG:H	1.56	0.46
35:DA:245:G:H5'	46:DP:70:GLN:H	1.80	0.46
47:DQ:127:ILE:HG22	47:DQ:128:LYS:N	2.21	0.46
47:DQ:86:GLY:C	47:DQ:88:GLY:H	2.20	0.46
50:DT:105:LEU:O	50:DT:113:LYS:NZ	2.49	0.46
45:DO:71:ARG:HH12	50:DT:74:ARG:HH22	1.63	0.46
51:DU:78:THR:C	51:DU:80:ILE:H	2.19	0.46
52:DV:3:ALA:HB3	52:DV:14:VAL:CB	2.45	0.46
28:D2:26:ARG:CD	54:DX:5:TYR:HB3	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:166:SER:OG	56:DZ:169:GLU:HB3	2.15	0.46
1:AA:1015:A:H2'	1:AA:1016:A:H8	1.81	0.45
1:AA:1062:U:H2'	1:AA:1063:C:C5	2.51	0.45
1:AA:1223:C:OP2	1:AA:1224:G:H8	1.98	0.45
1:AA:1295:G:O2'	13:AM:14:ARG:NH1	2.49	0.45
1:AA:1357:A:N6	1:AA:1363(A):A:H2	2.14	0.45
1:AA:927:G:H4'	1:AA:1503:A:N7	2.31	0.45
1:AA:27:G:H2'	1:AA:28:G:H8	1.81	0.45
1:AA:36:C:O2'	1:AA:37:U:H5'	2.15	0.45
1:AA:551:U:C2	1:AA:552:U:C5	3.04	0.45
1:AA:573:A:C2	1:AA:574:A:C2	3.05	0.45
1:AA:677:U:H3	1:AA:713:G:N2	2.12	0.45
4:AD:101:LEU:O	4:AD:104:VAL:HB	2.17	0.45
4:AD:13:ARG:O	4:AD:14:ARG:C	2.54	0.45
1:AA:409:G:H5'	4:AD:25:ARG:HB2	1.97	0.45
7:AG:133:GLY:HA2	7:AG:136:LYS:CG	2.46	0.45
7:AG:149:ARG:HD3	11:AK:59:TYR:CE1	2.51	0.45
7:AG:15:ASP:O	7:AG:19:GLY:HA2	2.16	0.45
7:AG:66:VAL:C	7:AG:68:ASN:H	2.18	0.45
8:AH:85:ARG:HA	8:AH:135:CYS:HB3	1.97	0.45
1:AA:1351:U:O4	9:AI:118:LYS:HE2	2.16	0.45
9:AI:47:LEU:CB	9:AI:50:LEU:HD12	2.46	0.45
9:AI:5:TYR:OH	9:AI:7:THR:HG23	2.16	0.45
9:AI:85:LEU:HD12	9:AI:85:LEU:C	2.36	0.45
10:AJ:4:ILE:HG23	10:AJ:98:ILE:CG2	2.41	0.45
16:AP:48:TRP:O	16:AP:49:LEU:C	2.52	0.45
16:AP:67:THR:HG21	16:AP:69:THR:HG23	1.98	0.45
17:AQ:67:LYS:C	17:AQ:70:ARG:HH12	2.18	0.45
25:AY:184:LEU:CD2	25:AY:185:GLY:N	2.79	0.45
25:AY:66:LEU:N	25:AY:66:LEU:HD12	2.31	0.45
26:B0:38:VAL:CG2	26:B0:59:LEU:HB2	2.46	0.45
35:BA:1215:G:H2'	35:BA:1216:G:O4'	2.16	0.45
35:BA:1341:U:H2'	35:BA:1397:U:O2	2.15	0.45
35:BA:1425:G:H2'	35:BA:1426:G:C8	2.51	0.45
35:BA:1711:C:H2'	35:BA:1712:C:C6	2.50	0.45
35:BA:729:G:C4	35:BA:1775:U:C2	3.03	0.45
35:BA:1935:G:C3'	35:BA:1962:C:H42	2.28	0.45
35:BA:1986:A:H3'	35:BA:1987:G:C5'	2.37	0.45
35:BA:2057:A:O2'	35:BA:2058:A:H5'	2.15	0.45
35:BA:2101:G:C6	35:BA:2102:U:C6	3.04	0.45
35:BA:2295:C:O2'	35:BA:2296:U:H5'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2523:G:H5'	35:BA:2523:G:H8	1.80	0.45
35:BA:2584:U:O4'	35:BA:2584:U:O2	2.34	0.45
35:BA:2650:U:H2'	35:BA:2651:C:H6	1.80	0.45
35:BA:2764:A:H2'	35:BA:2766:G:C8	2.51	0.45
35:BA:2793:G:C2	35:BA:2794:C:N3	2.85	0.45
35:BA:524:U:H4'	35:BA:555:U:H4'	1.97	0.45
35:BA:947:G:H2'	35:BA:948:G:H8	1.81	0.45
38:BD:49:ILE:O	38:BD:49:ILE:HG13	2.15	0.45
39:BE:48:GLN:CG	39:BE:78:LEU:HD12	2.45	0.45
41:BG:52:ILE:O	41:BG:54:GLU:HG3	2.15	0.45
35:BA:2312:U:O3'	41:BG:71:THR:HG21	2.15	0.45
35:BA:2745:C:H1'	42:BH:143:GLN:HG2	1.97	0.45
43:BI:2:LYS:O	43:BI:39:ALA:HB2	2.16	0.45
43:BI:62:LYS:C	43:BI:62:LYS:HD3	2.37	0.45
45:BO:1:MET:HE2	45:BO:1:MET:N	2.30	0.45
46:BP:56:SER:C	46:BP:58:THR:N	2.69	0.45
47:BQ:141:GLN:OE1	56:BZ:89:PHE:HB3	2.16	0.45
48:BR:74:LYS:O	48:BR:75:LEU:C	2.55	0.45
49:BS:74:ALA:HB2	49:BS:101:LEU:HD11	1.98	0.45
49:BS:15:ARG:HA	49:BS:17:ARG:HG2	1.98	0.45
50:BT:70:VAL:HG12	50:BT:71:GLY:O	2.16	0.45
50:BT:32:TYR:HD2	50:BT:81:PRO:CB	2.29	0.45
51:BU:111:GLU:O	51:BU:115:ALA:CB	2.64	0.45
52:BV:39:LEU:HD11	52:BV:53:GLU:CA	2.45	0.45
52:BV:72:VAL:CG1	52:BV:73:SER:H	2.21	0.45
53:BW:10:VAL:CG2	53:BW:101:SER:O	2.65	0.45
54:BX:59:VAL:HG22	54:BX:74:PRO:O	2.16	0.45
55:BY:37:VAL:CG2	55:BY:38:ILE:H	2.04	0.45
56:BZ:108:PRO:HB2	56:BZ:144:LEU:O	2.16	0.45
56:BZ:54:HIS:HB3	56:BZ:101:PRO:CD	2.46	0.45
56:BZ:79:ARG:O	56:BZ:79:ARG:HG2	2.15	0.45
1:CA:1098:C:C2	1:CA:1099:G:C8	3.04	0.45
1:CA:1118:C:O5'	9:CI:104:ARG:HG3	2.16	0.45
1:CA:1161:C:H2'	1:CA:1162:C:H6	1.81	0.45
1:CA:1262:C:C2	1:CA:1263:C:C5	3.05	0.45
1:CA:1462:G:O2'	1:CA:1463:C:H5'	2.16	0.45
1:CA:1493:A:H5''	1:CA:1494:G:OP1	2.15	0.45
1:CA:193:C:H2'	1:CA:194:C:C6	2.51	0.45
1:CA:193:C:O2'	1:CA:194:C:H5'	2.16	0.45
1:CA:321:A:N7	1:CA:328:C:O2	2.49	0.45
1:CA:445:G:C6	1:CA:490:G:O6	2.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:799:G:C2'	1:CA:800:G:H5'	2.46	0.45
2:CB:72:GLY:HA2	2:CB:165:VAL:HG22	1.98	0.45
2:CB:74:LYS:O	2:CB:75:LYS:C	2.55	0.45
2:CB:77:ALA:C	2:CB:80:ILE:HG23	2.37	0.45
4:CD:25:ARG:C	4:CD:27:TYR:N	2.69	0.45
5:CE:144:THR:O	5:CE:145:LYS:C	2.54	0.45
6:CF:9:VAL:HG12	6:CF:10:LEU:N	2.30	0.45
7:CG:150:ALA:C	7:CG:152:ALA:N	2.68	0.45
8:CH:35:ILE:HG22	8:CH:111:ILE:HD13	1.99	0.45
9:CI:18:PHE:HD1	9:CI:62:TYR:CD2	2.34	0.45
9:CI:48:GLU:N	9:CI:49:PRO:CD	2.79	0.45
1:CA:972:C:C4'	10:CJ:57:LYS:HG3	2.41	0.45
11:CK:86:GLY:N	11:CK:112:THR:HG23	2.29	0.45
13:CM:116:THR:HG22	13:CM:117:VAL:H	1.79	0.45
13:CM:10:PRO:HB2	13:CM:18:ALA:HB1	1.98	0.45
14:CN:41:ARG:HG3	14:CN:42:ILE:N	2.30	0.45
15:CO:41:GLU:O	15:CO:44:LYS:HB3	2.16	0.45
16:CP:23:ASP:OD1	16:CP:24:ALA:N	2.49	0.45
17:CQ:67:LYS:O	17:CQ:69:LYS:N	2.49	0.45
20:CT:74:LYS:CG	20:CT:75:ASN:H	2.28	0.45
25:CY:30:THR:O	25:CY:32:ARG:N	2.49	0.45
27:D1:64:ALA:O	27:D1:67:ILE:CG1	2.57	0.45
31:D5:49:CYS:O	31:D5:57:VAL:HG22	2.16	0.45
32:D6:26:ASN:O	32:D6:27:LYS:HD3	2.15	0.45
35:DA:585:G:C5	35:DA:1251:C:C4	3.04	0.45
35:DA:1300:U:O2'	35:DA:1301:A:P	2.74	0.45
35:DA:1453:U:H4'	35:DA:1455:G:OP1	2.16	0.45
35:DA:1568:G:P	38:DD:63:ARG:HH22	2.38	0.45
35:DA:1647:G:OP2	35:DA:1647:G:H3'	2.17	0.45
35:DA:1775:U:H2'	35:DA:1776:G:O5'	2.16	0.45
35:DA:1806:C:C5	35:DA:1807:G:N7	2.84	0.45
35:DA:858:U:O2	35:DA:2268:A:H2'	2.16	0.45
35:DA:271(W):G:H5'	35:DA:271(X):G:OP2	2.16	0.45
35:DA:2870:C:H2'	35:DA:2871:C:C5'	2.46	0.45
35:DA:371:A:C4	35:DA:373:U:C4	3.04	0.45
35:DA:614:U:O2	35:DA:614:U:O4'	2.31	0.45
35:DA:957:A:OP1	35:DA:957:A:H8	1.99	0.45
38:DD:131:LEU:H	38:DD:131:LEU:HD12	1.80	0.45
40:DF:55:GLY:O	40:DF:56:GLU:C	2.53	0.45
41:DG:113:ARG:C	41:DG:114:ILE:HG12	2.36	0.45
41:DG:38:VAL:HG13	41:DG:92:VAL:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:13:LYS:CE	42:DH:13:LYS:HA	2.36	0.45
42:DH:35:VAL:O	42:DH:37:VAL:HG23	2.16	0.45
42:DH:61:HIS:O	42:DH:65:HIS:HB3	2.17	0.45
42:DH:84:SER:O	42:DH:133:VAL:O	2.34	0.45
44:DN:122:VAL:HG12	44:DN:123:TYR:N	2.29	0.45
44:DN:42:TRP:HE3	44:DN:48:MET:SD	2.39	0.45
45:DO:10:VAL:HB	45:DO:12:ASP:OD2	2.15	0.45
45:DO:77:ILE:HD13	50:DT:74:ARG:HG3	1.99	0.45
46:DP:126:VAL:HG22	46:DP:145:PRO:CB	2.46	0.45
47:DQ:55:VAL:CG2	47:DQ:56:ARG:N	2.79	0.45
47:DQ:70:PRO:HA	47:DQ:95:ALA:HB2	1.98	0.45
48:DR:14:SER:O	48:DR:16:HIS:N	2.49	0.45
35:DA:1278:A:C5'	48:DR:36:THR:HG22	2.45	0.45
50:DT:102:ILE:O	50:DT:103:ARG:C	2.54	0.45
50:DT:52:ILE:HA	50:DT:61:PHE:HA	1.98	0.45
52:DV:22:VAL:CG2	52:DV:96:ILE:HB	2.45	0.45
52:DV:43:GLU:HA	52:DV:47:VAL:N	2.31	0.45
52:DV:83:ARG:CG	52:DV:83:ARG:NH1	2.68	0.45
53:DW:10:VAL:HB	53:DW:101:SER:O	2.17	0.45
54:DX:72:LYS:O	54:DX:73:ARG:HB3	2.16	0.45
55:DY:28:LYS:HZ1	55:DY:37:VAL:HA	1.74	0.45
55:DY:60:PHE:O	55:DY:62:GLU:OE2	2.34	0.45
1:AA:1279:A:N3	1:AA:1279:A:H2'	2.30	0.45
1:AA:1412:C:H2'	1:AA:1413:A:H8	1.70	0.45
1:AA:1488:G:H2'	1:AA:1489:G:C8	2.51	0.45
1:AA:16:A:N1	1:AA:919:A:C2	2.81	0.45
1:AA:241:C:C1'	1:AA:286:G:N2	2.79	0.45
1:AA:522:C:H1'	1:AA:536:C:H5'	1.97	0.45
1:AA:505:G:C6	1:AA:535:A:C2	3.04	0.45
1:AA:630:G:H2'	1:AA:631:G:C5'	2.45	0.45
1:AA:959:A:H3'	1:AA:960:U:H5''	1.99	0.45
1:AA:993:G:H22	1:AA:1046:A:H1'	1.82	0.45
2:AB:51:LEU:O	2:AB:55:PHE:HD2	1.98	0.45
3:AC:207:VAL:HG12	3:AC:207:VAL:O	2.15	0.45
5:AE:55:VAL:O	5:AE:58:ALA:HB3	2.16	0.45
7:AG:104:LEU:HD22	7:AG:104:LEU:H	1.80	0.45
7:AG:148:ASN:O	7:AG:150:ALA:N	2.50	0.45
8:AH:114:THR:HG21	8:AH:119:LEU:HD21	1.97	0.45
8:AH:11:THR:CA	8:AH:14:ARG:HH12	2.29	0.45
10:AJ:28:ARG:NH2	10:AJ:34:VAL:O	2.50	0.45
1:AA:1150:U:O3'	10:AJ:41:PRO:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AK:21:ILE:HG13	11:AK:30:VAL:HG12	1.98	0.45
11:AK:23:ALA:O	11:AK:87:THR:N	2.49	0.45
14:AN:27:CYS:O	14:AN:27:CYS:SG	2.74	0.45
10:AJ:65:LEU:HD12	14:AN:55:GLY:HA3	1.97	0.45
16:AP:39:TYR:C	16:AP:39:TYR:CD1	2.90	0.45
20:AT:93:GLU:OE1	20:AT:93:GLU:N	2.49	0.45
28:B2:26:ARG:O	28:B2:29:LYS:N	2.48	0.45
29:B3:19:GLN:O	29:B3:21:ALA:N	2.50	0.45
32:B6:27:LYS:HE2	35:BA:2285:C:H5	1.81	0.45
33:B7:23:ARG:O	33:B7:28:ARG:NH1	2.48	0.45
35:BA:105:C:H5'	35:BA:106:C:OP2	2.16	0.45
35:BA:9:U:O2'	35:BA:10:G:O5'	2.33	0.45
35:BA:1186:G:O2'	35:BA:1187:G:H5'	2.16	0.45
35:BA:1315:C:H42	35:BA:1337:G:H1	1.63	0.45
35:BA:225:A:O2'	35:BA:257:A:H4'	2.16	0.45
35:BA:2606:C:C2'	35:BA:2607:G:H5'	2.46	0.45
35:BA:2695:C:H2'	35:BA:2696:U:C6	2.50	0.45
35:BA:60:G:C5	35:BA:63:U:C4	3.04	0.45
35:BA:76:C:H2'	35:BA:77:C:C6	2.51	0.45
35:BA:806:C:O2'	35:BA:807:U:H5'	2.16	0.45
35:BA:966:G:H2'	35:BA:967:C:H6	1.80	0.45
38:BD:133:LEU:HB2	38:BD:187:GLY:HA2	1.98	0.45
39:BE:120:TRP:CD2	39:BE:155:LYS:HD3	2.51	0.45
39:BE:116:VAL:CG2	39:BE:122:PHE:CG	2.99	0.45
39:BE:88:GLY:O	39:BE:89:ASP:CB	2.64	0.45
40:BF:24:LEU:O	40:BF:25:PRO:C	2.53	0.45
40:BF:25:PRO:HB3	40:BF:119:ARG:HD3	1.98	0.45
40:BF:28:ILE:CD1	40:BF:28:ILE:N	2.78	0.45
45:BO:60:ALA:HA	45:BO:87:ILE:H	1.82	0.45
46:BP:25:SER:OG	46:BP:26:GLY:N	2.47	0.45
46:BP:70:GLN:CG	46:BP:71:VAL:N	2.77	0.45
49:BS:26:LEU:CD2	49:BS:28:VAL:HG22	2.46	0.45
51:BU:111:GLU:C	51:BU:113:ALA:N	2.68	0.45
52:BV:3:ALA:HB3	52:BV:14:VAL:CB	2.42	0.45
52:BV:75:PHE:CE1	52:BV:89:GLN:HB2	2.51	0.45
54:BX:12:VAL:HG13	54:BX:17:ALA:HB2	1.96	0.45
54:BX:62:LYS:CD	54:BX:68:ARG:HD2	2.47	0.45
55:BY:34:LYS:O	55:BY:35:TYR:HB3	2.17	0.45
56:BZ:71:VAL:HG13	56:BZ:86:VAL:HG13	1.97	0.45
1:CA:1006:C:H2'	1:CA:1007:C:C6	2.51	0.45
1:CA:1054:C:O2'	1:CA:1055:A:H5''	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1157:A:H1'	1:CA:1181:G:N2	2.32	0.45
1:CA:137:C:H2'	1:CA:138:G:C8	2.52	0.45
1:CA:1486:G:C6	1:CA:1487:G:C6	3.05	0.45
1:CA:243:A:C2	1:CA:245:C:C2	3.05	0.45
1:CA:261:U:O2	1:CA:263:A:C8	2.68	0.45
1:CA:726:C:H2'	1:CA:727:G:H8	1.80	0.45
1:CA:76:C:N4	1:CA:93:G:H1	2.14	0.45
1:CA:902:G:O2'	1:CA:903:G:H5'	2.16	0.45
1:CA:929:G:O2'	1:CA:930:C:H5'	2.16	0.45
2:CB:221:LEU:O	2:CB:221:LEU:HD22	2.15	0.45
5:CE:60:TYR:CE2	5:CE:64:ARG:NH2	2.84	0.45
1:CA:1351:U:O4	9:CI:118:LYS:HE2	2.16	0.45
9:CI:17:VAL:HG21	9:CI:81:ILE:HD13	1.96	0.45
9:CI:9:ARG:HA	9:CI:14:VAL:HA	1.99	0.45
11:CK:33:THR:HG22	11:CK:39:PRO:HA	1.98	0.45
12:CL:37:CYS:O	12:CL:79:GLU:O	2.34	0.45
12:CL:95:GLY:O	12:CL:97:ARG:N	2.49	0.45
13:CM:91:ARG:HB2	13:CM:98:VAL:HG22	1.98	0.45
14:CN:13:THR:N	14:CN:14:PRO:CD	2.80	0.45
1:CA:617:G:H4'	16:CP:44:THR:O	2.16	0.45
18:CR:36:ASN:HB2	18:CR:39:VAL:CG2	2.46	0.45
20:CT:25:ARG:CZ	20:CT:25:ARG:HB2	2.46	0.45
25:CY:108:GLU:O	25:CY:109:GLU:C	2.53	0.45
25:CY:134:ARG:HG2	25:CY:138:ASP:OD1	2.16	0.45
25:CY:46:TYR:C	25:CY:48:ALA:H	2.17	0.45
25:CY:43:VAL:HB	25:CY:50:VAL:HG23	1.98	0.45
25:CY:83:ILE:HG22	25:CY:84:ARG:N	2.31	0.45
26:D0:11:ARG:O	26:D0:12:ASN:OD1	2.33	0.45
27:D1:87:PRO:CB	27:D1:91:LYS:HZ2	2.29	0.45
34:D8:30:ARG:NE	46:DP:62:LEU:HB2	2.29	0.45
35:DA:120:U:O4	35:DA:177:G:C8	2.69	0.45
35:DA:1866:C:H2'	35:DA:1876:A:O4'	2.15	0.45
35:DA:1937:A:O2'	35:DA:1938:A:OP1	2.26	0.45
35:DA:2034:U:H2'	35:DA:2035:G:H5'	1.98	0.45
35:DA:2352:A:C2'	35:DA:2353:G:H5'	2.46	0.45
35:DA:2704:C:C4	35:DA:2705:A:N7	2.84	0.45
35:DA:563:G:OP2	35:DA:572:A:H5'	2.15	0.45
35:DA:580:C:H2'	35:DA:581:C:H6	1.80	0.45
35:DA:71:A:C8	35:DA:71:A:H5'	2.51	0.45
35:DA:826:U:H3'	35:DA:828:U:C6	2.51	0.45
35:DA:852:G:C2'	35:DA:853:G:H5'	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:109:ASP:CB	38:DD:195:ALA:HB3	2.45	0.45
38:DD:61:LEU:HD12	38:DD:62:TYR:H	1.80	0.45
38:DD:25:THR:HG21	38:DD:82:ILE:H	1.81	0.45
39:DE:61:ARG:CG	39:DE:62:PRO:HD3	2.44	0.45
41:DG:21:ARG:HH11	41:DG:22:ARG:HB2	1.81	0.45
43:DI:110:ASP:O	43:DI:114:LEU:HG	2.17	0.45
44:DN:38:HIS:CG	44:DN:39:ARG:H	2.34	0.45
45:DO:14:THR:HG21	45:DO:86:ILE:CG1	2.46	0.45
52:DV:17:GLY:HA2	52:DV:98:GLU:O	2.16	0.45
52:DV:75:PHE:HB2	52:DV:87:HIS:HB3	1.97	0.45
53:DW:10:VAL:O	53:DW:11:ARG:CB	2.64	0.45
53:DW:92:ARG:HG2	53:DW:92:ARG:HH11	1.80	0.45
1:AA:1168:A:H2'	1:AA:1169:A:H8	1.79	0.45
1:AA:1226:C:C5'	19:AS:80:TYR:HE2	2.29	0.45
1:AA:122:G:H2'	1:AA:123:C:C6	2.52	0.45
1:AA:1375:A:C4	1:AA:1376:U:C5	3.04	0.45
1:AA:302:G:C4	1:AA:303:A:C8	3.04	0.45
1:AA:486:U:H2'	1:AA:487:A:H8	1.81	0.45
1:AA:445:G:C6	1:AA:490:G:O6	2.70	0.45
1:AA:62:U:H5''	1:AA:385:C:O2	2.16	0.45
1:AA:659:U:H2'	1:AA:660:G:C8	2.50	0.45
1:AA:908:A:C2	1:AA:909:A:N7	2.84	0.45
2:AB:178:ARG:CG	2:AB:178:ARG:HH11	2.29	0.45
2:AB:46:LYS:O	2:AB:47:THR:C	2.54	0.45
3:AC:95:THR:HG22	3:AC:97:LYS:HB2	1.96	0.45
4:AD:194:LEU:N	4:AD:194:LEU:CD2	2.74	0.45
4:AD:88:VAL:HG12	4:AD:88:VAL:O	2.16	0.45
6:AF:67:MET:CE	6:AF:72:VAL:H	2.29	0.45
7:AG:30:ILE:HD13	7:AG:105:VAL:HG13	1.98	0.45
7:AG:57:GLU:O	7:AG:59:LEU:N	2.46	0.45
7:AG:86:GLN:NE2	23:AW:32:G:N2	2.63	0.45
9:AI:28:VAL:HG13	9:AI:63:ILE:C	2.36	0.45
11:AK:126:ARG:O	11:AK:127:LYS:C	2.54	0.45
12:AL:18:VAL:O	12:AL:18:VAL:HG23	2.16	0.45
12:AL:84:LEU:HB3	12:AL:101:VAL:HB	1.98	0.45
20:AT:27:LYS:HD3	20:AT:27:LYS:C	2.36	0.45
25:AY:111:ARG:O	25:AY:115:VAL:HG23	2.16	0.45
25:AY:15:GLN:HA	25:AY:168:PHE:CZ	2.51	0.45
25:AY:64:ARG:O	25:AY:65:THR:OG1	2.31	0.45
25:AY:74:ASN:HA	25:AY:77:LYS:HG2	1.98	0.45
27:B1:16:ASN:HB3	27:B1:46:LEU:CG	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:81:LYS:O	27:B1:82:LEU:C	2.55	0.45
28:B2:49:LYS:O	28:B2:52:ASP:HB3	2.15	0.45
31:B5:30:LEU:HD23	31:B5:41:PRO:HB3	1.97	0.45
35:BA:1277:G:H2'	35:BA:1278:A:H8	1.81	0.45
35:BA:1331:A:H2'	35:BA:1333:C:C5	2.51	0.45
35:BA:154(A):C:O4'	35:BA:154(A):C:O2	2.34	0.45
35:BA:195:A:H5''	35:BA:196:A:OP2	2.16	0.45
35:BA:2283:C:C5	35:BA:2389:G:H2'	2.51	0.45
35:BA:2663:G:H2'	35:BA:2664:G:O4'	2.16	0.45
35:BA:2775:A:O2'	35:BA:2776:A:C5'	2.65	0.45
35:BA:601:C:O2'	35:BA:605:C:H5''	2.16	0.45
35:BA:60:G:N2	35:BA:74:A:H2'	2.31	0.45
35:BA:609:A:H2'	35:BA:610:G:O4'	2.17	0.45
35:BA:914:C:C2'	35:BA:915:C:H5'	2.41	0.45
38:BD:109:ASP:CB	38:BD:195:ALA:HB3	2.46	0.45
39:BE:1:MET:H3	39:BE:84:PHE:HB2	1.82	0.45
40:BF:169:ASN:ND2	40:BF:169:ASN:O	2.49	0.45
40:BF:45:ARG:NH1	40:BF:97:TYR:CD2	2.84	0.45
41:BG:149:VAL:O	41:BG:149:VAL:HG23	2.17	0.45
41:BG:171:ALA:O	41:BG:173:LEU:N	2.49	0.45
44:BN:38:HIS:CG	44:BN:39:ARG:N	2.84	0.45
47:BQ:70:PRO:HA	47:BQ:94:VAL:O	2.15	0.45
53:BW:8:ARG:NH1	53:BW:8:ARG:HG3	2.31	0.45
56:BZ:134:PRO:O	56:BZ:135:GLU:C	2.54	0.45
1:CA:1011:G:O2'	1:CA:1012:U:H5'	2.17	0.45
1:CA:1030(D):A:C2'	1:CA:1031:G:H5'	2.45	0.45
1:CA:1153:C:H2'	1:CA:1154:G:O5'	2.17	0.45
1:CA:311:C:O2'	1:CA:312:C:H5'	2.15	0.45
1:CA:502:G:C6	1:CA:544:G:C6	3.04	0.45
1:CA:650:G:O2'	1:CA:651:C:H5'	2.16	0.45
1:CA:780:A:H2	1:CA:803:G:C6	2.32	0.45
1:CA:894:G:H2'	1:CA:895:G:C8	2.51	0.45
2:CB:100:GLY:O	2:CB:104:ASN:N	2.49	0.45
3:CC:92:ALA:C	3:CC:94:LEU:N	2.70	0.45
4:CD:22:LYS:HB3	4:CD:22:LYS:HZ2	1.81	0.45
4:CD:32:ALA:O	4:CD:36:ARG:N	2.49	0.45
5:CE:101:ILE:O	5:CE:120:THR:OG1	2.34	0.45
6:CF:42:GLU:O	6:CF:44:GLY:N	2.50	0.45
7:CG:6:ARG:O	7:CG:7:ALA:C	2.55	0.45
8:CH:14:ARG:HB3	8:CH:14:ARG:HH11	1.80	0.45
8:CH:87:SER:OG	8:CH:92:ARG:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CI:92:TYR:HB3	9:CI:95:LYS:HD2	1.99	0.45
1:CA:568:G:O6	12:CL:5:PRO:HD3	2.16	0.45
13:CM:54:VAL:O	13:CM:58:GLU:HG2	2.16	0.45
16:CP:49:LEU:HD11	16:CP:51:VAL:CG2	2.38	0.45
16:CP:80:PHE:N	16:CP:80:PHE:CD1	2.85	0.45
17:CQ:57:VAL:HA	17:CQ:77:VAL:HG23	1.99	0.45
18:CR:31:LEU:HG	18:CR:65:ILE:HD13	1.99	0.45
25:CY:31:GLY:O	25:CY:32:ARG:HB2	2.16	0.45
25:CY:45:TYR:HD2	25:CY:78:ALA:HB2	1.80	0.45
25:CY:83:ILE:O	25:CY:85:ASP:N	2.50	0.45
27:D1:23:LYS:O	27:D1:37:ILE:CG1	2.65	0.45
27:D1:75:GLU:HB2	27:D1:76:ARG:HH21	1.76	0.45
28:D2:21:LEU:HD23	28:D2:24:LEU:HD22	1.98	0.45
29:D3:38:GLU:CD	29:D3:38:GLU:H	2.20	0.45
29:D3:40:THR:OG1	29:D3:41:PRO:HD2	2.16	0.45
35:DA:1221(A):C:C2'	35:DA:1222:C:H5'	2.45	0.45
35:DA:1247:A:OP1	40:DF:95:ARG:NH2	2.49	0.45
35:DA:1341:U:H2'	35:DA:1397:U:O2	2.16	0.45
35:DA:1497:U:H3	35:DA:1578:U:P	2.39	0.45
35:DA:151:C:H2'	35:DA:152:G:H8	1.81	0.45
35:DA:1813:G:H1'	38:DD:50:THR:OG1	2.16	0.45
35:DA:1820:U:H4'	35:DA:1821:A:OP2	2.17	0.45
35:DA:1945:G:C6	35:DA:1946:U:O4	2.70	0.45
35:DA:2171:A:HO2'	35:DA:2172:U:H6	1.60	0.45
35:DA:2428:G:H5''	35:DA:2429:G:O5'	2.16	0.45
35:DA:2474:C:O2	35:DA:2474:C:H2'	2.14	0.45
35:DA:2523:G:H5'	35:DA:2523:G:H8	1.81	0.45
35:DA:2663:G:H2'	35:DA:2664:G:O4'	2.16	0.45
35:DA:2000:G:O2'	35:DA:2689:U:H5	1.96	0.45
35:DA:290:G:O2'	35:DA:291:C:H5'	2.16	0.45
35:DA:395:U:H1'	35:DA:396:G:N7	2.32	0.45
35:DA:64:A:C2	35:DA:65:C:C2	3.03	0.45
35:DA:70:G:H21	35:DA:71:A:N6	2.15	0.45
35:DA:823:G:H2'	35:DA:824:A:H8	1.81	0.45
35:DA:897:C:O2'	35:DA:898:C:H5'	2.16	0.45
37:DC:77:ILE:HG13	37:DC:96:GLY:O	2.15	0.45
37:DC:86:ALA:HA	37:DC:89:ALA:HB2	1.99	0.45
41:DG:39:ILE:CB	41:DG:157:ILE:HG22	2.45	0.45
42:DH:146:ALA:O	42:DH:148:ILE:N	2.50	0.45
42:DH:39:PRO:O	42:DH:42:ARG:HG3	2.17	0.45
42:DH:43:VAL:CG2	42:DH:43:VAL:O	2.61	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:87:ILE:HD13	45:DO:87:ILE:HA	1.79	0.45
48:DR:4:LEU:HD22	48:DR:4:LEU:O	2.17	0.45
54:DX:36:LYS:O	54:DX:37:THR:C	2.54	0.45
55:DY:15:VAL:O	55:DY:16:ALA:HB2	2.17	0.45
55:DY:19:LYS:HG2	55:DY:19:LYS:O	2.15	0.45
1:AA:1019:C:H2'	1:AA:1020:U:O4'	2.17	0.45
1:AA:1164:G:H2'	1:AA:1165:C:H5'	1.97	0.45
1:AA:1289:A:H3'	1:AA:1290:G:H8	1.81	0.45
1:AA:137:C:H2'	1:AA:138:G:C8	2.51	0.45
1:AA:66:G:C4'	1:AA:173:U:C4	3.00	0.45
1:AA:557:G:H2'	1:AA:558:G:C8	2.51	0.45
1:AA:695:A:O2'	1:AA:696:A:H5'	2.16	0.45
1:AA:707:C:H2'	1:AA:707:C:O2	2.17	0.45
1:AA:918:A:H2'	1:AA:919:A:C8	2.51	0.45
1:AA:972:C:H4'	10:AJ:57:LYS:HG2	1.97	0.45
2:AB:195:ASP:C	2:AB:197:VAL:H	2.20	0.45
2:AB:212:GLN:NE2	2:AB:235:SER:HB3	2.31	0.45
3:AC:134:ILE:HG22	3:AC:168:ALA:CB	2.47	0.45
3:AC:173:VAL:N	3:AC:174:PRO:CD	2.78	0.45
4:AD:102:ASP:HA	4:AD:121:VAL:HG21	1.97	0.45
4:AD:18:LYS:NZ	4:AD:31:CYS:CB	2.77	0.45
4:AD:34:GLU:O	4:AD:35:ARG:HG2	2.16	0.45
7:AG:15:ASP:HB3	7:AG:20:ASP:H	1.81	0.45
8:AH:5:PRO:C	8:AH:8:ASP:HB3	2.37	0.45
11:AK:48:ILE:C	11:AK:50:TYR:H	2.20	0.45
14:AN:13:THR:N	14:AN:14:PRO:CD	2.79	0.45
16:AP:80:PHE:CD1	16:AP:80:PHE:N	2.83	0.45
18:AR:30:ASP:O	18:AR:32:ARG:N	2.48	0.45
19:AS:52:TYR:CG	19:AS:53:ASN:N	2.84	0.45
21:AU:12:LYS:O	21:AU:22:ARG:NH1	2.49	0.45
1:AA:1328:C:P	21:AU:21:TYR:OH	2.74	0.45
25:AY:171:LYS:O	25:AY:175:LEU:HB2	2.15	0.45
25:AY:175:LEU:O	25:AY:176:ALA:C	2.53	0.45
26:B0:11:ARG:O	26:B0:12:ASN:OD1	2.34	0.45
27:B1:75:GLU:O	27:B1:76:ARG:NE	2.50	0.45
29:B3:12:PRO:O	29:B3:14:GLY:N	2.50	0.45
29:B3:40:THR:O	29:B3:43:ILE:N	2.50	0.45
35:BA:1112:G:O2'	35:BA:1113:U:H5''	2.17	0.45
35:BA:1263:U:C5	35:BA:1264:G:C6	3.05	0.45
35:BA:132:G:H2'	35:BA:133:C:C6	2.51	0.45
35:BA:1861:G:H2'	35:BA:1862:G:H8	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:186:G:C2'	35:BA:187:G:H5'	2.47	0.45
35:BA:1998:G:H4'	35:BA:2724:C:H4'	1.99	0.45
35:BA:2023:G:C2	35:BA:2024:G:C5	3.05	0.45
35:BA:2035:G:H4'	35:BA:2036:C:OP2	2.16	0.45
35:BA:2319:G:C5	35:BA:2320:A:N6	2.85	0.45
35:BA:272(E):G:C2	35:BA:364:C:N3	2.85	0.45
35:BA:2832:U:O4	35:BA:2883:A:H5''	2.16	0.45
35:BA:360:G:H2'	35:BA:361:G:C8	2.50	0.45
35:BA:412:A:H3'	35:BA:413:C:C6	2.51	0.45
26:B0:77:ARG:NH2	35:BA:857:C:OP2	2.48	0.45
35:BA:910:A:N1	35:BA:2277:G:H1'	2.31	0.45
36:BB:100:A:C4	36:BB:101:G:C8	3.05	0.45
38:BD:66:ASP:OD2	38:BD:103:ARG:NH2	2.49	0.45
35:BA:574:C:N3	39:BE:145:LYS:HE2	2.32	0.45
39:BE:6:GLY:HA3	39:BE:27:LEU:O	2.16	0.45
40:BF:1:MET:O	40:BF:3:GLU:N	2.49	0.45
40:BF:45:ARG:HG3	40:BF:46:ARG:N	2.31	0.45
41:BG:53:LEU:CD2	41:BG:53:LEU:H	2.22	0.45
42:BH:121:ILE:CG2	42:BH:133:VAL:HG11	2.46	0.45
42:BH:37:VAL:HG12	42:BH:38:SER:N	2.31	0.45
42:BH:66:GLY:O	42:BH:67:LEU:C	2.54	0.45
44:BN:27:ALA:HB3	44:BN:106:MET:HE2	1.99	0.45
44:BN:36:GLY:HA3	44:BN:48:MET:CE	2.45	0.45
45:BO:37:ASP:O	45:BO:39:ILE:HG22	2.16	0.45
45:BO:67:LYS:O	45:BO:68:GLU:C	2.55	0.45
48:BR:32:GLY:O	48:BR:116:LEU:N	2.48	0.45
48:BR:73:VAL:HG23	48:BR:74:LYS:N	2.31	0.45
49:BS:86:ALA:C	49:BS:106:ARG:HG3	2.36	0.45
50:BT:122:ASP:O	50:BT:126:ALA:HB2	2.17	0.45
45:BO:104:ARG:NH1	50:BT:35:LYS:HB3	2.32	0.45
50:BT:35:LYS:CE	50:BT:41:ARG:HG3	2.44	0.45
52:BV:71:LEU:CD1	52:BV:72:VAL:N	2.79	0.45
35:BA:493:G:HO2'	53:BW:8:ARG:H	1.64	0.45
1:CA:1288:A:O5'	1:CA:1288:A:H8	1.99	0.45
1:CA:1473:A:O2'	35:DA:1702:G:H4'	2.17	0.45
1:CA:788:U:C2	1:CA:789:U:C6	3.04	0.45
1:CA:835:U:H3	1:CA:851:G:H1	1.65	0.45
1:CA:885:G:H1	1:CA:912:C:H42	1.64	0.45
2:CB:112:VAL:C	2:CB:115:LEU:HB3	2.32	0.45
3:CC:99:VAL:HG23	3:CC:99:VAL:O	2.15	0.45
4:CD:3:ARG:HG2	4:CD:3:ARG:HH21	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:76:ILE:CG2	5:CE:77:PRO:N	2.79	0.45
6:CF:83:ASP:C	6:CF:85:VAL:N	2.70	0.45
7:CG:26:PHE:CE1	7:CG:105:VAL:HG22	2.51	0.45
8:CH:53:VAL:O	8:CH:54:ASP:CB	2.64	0.45
16:CP:60:LEU:O	16:CP:62:VAL:N	2.49	0.45
1:CA:255:G:H5''	17:CQ:17:LYS:HB2	1.98	0.45
17:CQ:62:SER:OG	17:CQ:72:ARG:HG3	2.16	0.45
19:CS:11:VAL:HG22	19:CS:16:LEU:HD11	1.97	0.45
20:CT:100:ILE:H	20:CT:100:ILE:CD1	2.28	0.45
20:CT:24:LEU:O	20:CT:27:LYS:HB3	2.15	0.45
23:CW:52:C:N4	23:CW:53:G:C6	2.84	0.45
32:D6:52:VAL:HG12	32:D6:52:VAL:O	2.16	0.45
33:D7:18:PHE:O	33:D7:21:ARG:N	2.49	0.45
33:D7:5:TRP:HE1	33:D7:7:PRO:HG3	1.80	0.45
35:DA:1262:A:P	53:DW:99:ARG:HH12	2.40	0.45
35:DA:1264:G:C3'	35:DA:1265:A:H5''	2.46	0.45
27:D1:8:SER:HB3	35:DA:1364:G:C5'	2.46	0.45
35:DA:147:U:H2'	35:DA:148:C:C6	2.51	0.45
35:DA:1550:C:H2'	35:DA:1551:C:C6	2.41	0.45
35:DA:1566:A:C4	38:DD:214:TRP:CE3	3.05	0.45
35:DA:1608:A:H1'	35:DA:1610:A:OP2	2.17	0.45
35:DA:179:G:C4	35:DA:180:G:C8	3.03	0.45
35:DA:1794:U:HO2'	35:DA:1900:A:HO2'	1.61	0.45
35:DA:2225:A:H1'	35:DA:2226:C:OP2	2.17	0.45
35:DA:2308:G:N7	35:DA:2310:A:H5'	2.31	0.45
35:DA:2350:C:O2'	35:DA:2351:G:H5'	2.16	0.45
35:DA:2355:C:C4	35:DA:2356:C:C4	3.05	0.45
35:DA:2550:G:H2'	35:DA:2551:C:C6	2.51	0.45
35:DA:2801:A:O2'	35:DA:2895:U:C4'	2.65	0.45
35:DA:372:G:O2'	35:DA:373:U:P	2.75	0.45
35:DA:548:A:O2'	35:DA:549:G:OP1	2.26	0.45
35:DA:703:U:H2'	35:DA:704:G:C5'	2.43	0.45
36:DB:103:G:O2'	36:DB:104:U:H5'	2.17	0.45
38:DD:45:ASN:C	38:DD:46:GLN:OE1	2.55	0.45
38:DD:81:ALA:N	38:DD:94:LEU:CD1	2.79	0.45
38:DD:80:ALA:CB	38:DD:96:HIS:CE1	2.99	0.45
41:DG:96:ARG:O	41:DG:100:TRP:CD1	2.69	0.45
41:DG:102:PHE:CZ	41:DG:141:PHE:CE1	3.04	0.45
41:DG:61:ALA:O	41:DG:62:LEU:HD23	2.17	0.45
42:DH:149:ARG:CA	42:DH:162:ILE:HD11	2.45	0.45
42:DH:149:ARG:CB	42:DH:162:ILE:HD11	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:97:ARG:O	42:DH:98:LEU:HB2	2.15	0.45
44:DN:19:GLU:C	44:DN:21:LYS:H	2.20	0.45
46:DP:96:THR:HB	46:DP:126:VAL:HB	1.98	0.45
47:DQ:70:PRO:HA	47:DQ:94:VAL:O	2.16	0.45
48:DR:17:ARG:O	48:DR:18:LEU:C	2.53	0.45
48:DR:20:LEU:C	48:DR:20:LEU:CD1	2.75	0.45
45:DO:104:ARG:NH2	50:DT:33:LYS:HD2	2.31	0.45
51:DU:61:TRP:O	51:DU:62:ILE:C	2.55	0.45
52:DV:61:VAL:CG1	52:DV:62:LEU:N	2.79	0.45
52:DV:5:VAL:CG2	52:DV:6:LYS:N	2.80	0.45
53:DW:24:ILE:O	53:DW:71:VAL:HG11	2.16	0.45
55:DY:2:ARG:HG2	55:DY:2:ARG:NH1	2.32	0.45
56:DZ:110:GLY:O	56:DZ:111:VAL:CG1	2.64	0.45
56:DZ:45:ASP:O	56:DZ:47:VAL:N	2.49	0.45
56:DZ:76:LEU:HA	56:DZ:84:GLU:N	2.30	0.45
1:AA:1095:U:H2'	1:AA:1096:C:H6	1.82	0.45
1:AA:1412:C:C2'	1:AA:1413:A:C8	2.83	0.45
1:AA:1423:G:H2'	1:AA:1424:C:H6	1.81	0.45
1:AA:1443:G:C6	1:AA:1460:A:C2	3.04	0.45
1:AA:245:C:O2	1:AA:283:C:N3	2.49	0.45
1:AA:389:A:H2'	1:AA:390:C:H5'	1.99	0.45
1:AA:411:A:H2'	1:AA:412:A:H4'	1.99	0.45
1:AA:650:G:O2'	1:AA:651:C:H5'	2.16	0.45
1:AA:693:G:N2	23:AW:38:A:C2	2.77	0.45
1:AA:865:A:C2	1:AA:918:A:H4'	2.52	0.45
1:AA:866:C:C2	1:AA:867:G:H1'	2.51	0.45
1:AA:885:G:H2'	1:AA:886:G:H8	1.81	0.45
2:AB:115:LEU:HA	2:AB:145:LEU:HD11	1.99	0.45
2:AB:43:ASP:OD2	2:AB:46:LYS:N	2.44	0.45
3:AC:25:GLY:O	3:AC:28:GLN:N	2.49	0.45
3:AC:92:ALA:HB2	3:AC:99:VAL:HG22	1.97	0.45
7:AG:140:ASP:O	7:AG:143:ARG:HB2	2.17	0.45
8:AH:8:ASP:O	8:AH:12:ARG:HG3	2.16	0.45
9:AI:118:LYS:HB2	9:AI:121:ARG:HB2	1.97	0.45
9:AI:13:ALA:HB2	9:AI:68:GLY:HA3	1.98	0.45
10:AJ:8:LEU:CA	10:AJ:96:ILE:HG22	2.45	0.45
11:AK:23:ALA:CB	11:AK:91:ARG:HG2	2.45	0.45
13:AM:86:CYS:SG	13:AM:89:GLY:N	2.90	0.45
14:AN:47:LEU:HB3	14:AN:53:LEU:HG	1.98	0.45
1:AA:255:G:H5''	17:AQ:17:LYS:HB2	1.99	0.45
20:AT:58:LYS:O	20:AT:62:LEU:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:AV:41:C:H2'	22:AV:42:C:C6	2.51	0.45
25:AY:155:LYS:HA	25:AY:158:GLU:HB3	1.98	0.45
26:B0:25:ARG:HG3	26:B0:29:GLN:HE21	1.81	0.45
26:B0:42:GLY:HA3	35:BA:2331:G:O4'	2.16	0.45
27:B1:69:LYS:NZ	27:B1:95:LEU:HD12	2.31	0.45
35:BA:1048:A:N6	35:BA:1106:A:C8	2.84	0.45
35:BA:1204:A:N1	35:BA:1241:A:C2	2.85	0.45
35:BA:1810:A:H2'	35:BA:1811:G:O4'	2.16	0.45
35:BA:2003:G:C6	35:BA:2004:G:C5	3.05	0.45
35:BA:2322:A:H2'	35:BA:2323:G:O4'	2.16	0.45
35:BA:2383:G:H2'	35:BA:2384:G:C8	2.51	0.45
35:BA:2059:A:C5	35:BA:2503:A:C2	3.04	0.45
35:BA:2672:G:C3'	35:BA:2673:G:H5''	2.45	0.45
35:BA:2704:C:C4	35:BA:2705:A:N7	2.85	0.45
35:BA:290:G:N2	35:BA:291:C:H1'	2.32	0.45
35:BA:323:G:O2'	35:BA:1205:U:N3	2.43	0.45
35:BA:347:A:H2'	35:BA:348:G:H8	1.80	0.45
35:BA:366:C:C4	35:BA:404:C:H5	2.34	0.45
35:BA:532:A:H4'	35:BA:533:G:C1'	2.46	0.45
35:BA:618:C:H2'	35:BA:619:G:C8	2.52	0.45
35:BA:619:G:H5''	35:BA:620:G:OP2	2.17	0.45
35:BA:619:G:H3'	35:BA:620:G:N2	2.31	0.45
35:BA:607:U:H3	35:BA:621:A:H2	1.64	0.45
35:BA:660:G:H4'	40:BF:38:ARG:NH1	2.32	0.45
35:BA:675:A:C8	35:BA:804:A:C6	3.05	0.45
35:BA:916:G:H2'	35:BA:917:A:H5''	1.98	0.45
35:BA:977:G:C6	35:BA:987:G:C6	3.05	0.45
38:BD:133:LEU:HD12	38:BD:186:HIS:O	2.17	0.45
38:BD:245:PRO:O	38:BD:246:PRO:C	2.54	0.45
38:BD:45:ASN:C	38:BD:46:GLN:OE1	2.54	0.45
41:BG:46:ALA:HA	41:BG:51:ARG:HG3	1.96	0.45
43:BI:109:ILE:HD12	43:BI:111:PRO:HD3	1.98	0.45
43:BI:98:ALA:O	43:BI:109:ILE:HG21	2.17	0.45
46:BP:16:ARG:NH1	46:BP:18:ARG:CG	2.79	0.45
46:BP:48:PRO:CG	46:BP:49:ARG:N	2.79	0.45
46:BP:56:SER:O	46:BP:57:THR:CB	2.65	0.45
47:BQ:53:ALA:CA	47:BQ:56:ARG:HB3	2.47	0.45
47:BQ:68:ILE:H	47:BQ:68:ILE:CD1	2.21	0.45
47:BQ:76:LYS:N	47:BQ:88:GLY:HA3	2.31	0.45
50:BT:92:GLY:C	50:BT:94:ALA:H	2.18	0.45
52:BV:5:VAL:CG2	52:BV:6:LYS:N	2.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:9:TYR:CD2	53:BW:9:TYR:N	2.83	0.45
56:BZ:143:GLY:C	56:BZ:144:LEU:HD22	2.37	0.45
56:BZ:166:SER:CB	56:BZ:168:GLU:N	2.79	0.45
1:CA:1015:A:H8	1:CA:1015:A:O5'	1.99	0.45
1:CA:1040:U:H2'	1:CA:1041:A:H8	1.81	0.45
1:CA:1268:A:H2'	1:CA:1269:A:C8	2.52	0.45
1:CA:189(K):U:H2'	1:CA:189(L):G:C8	2.51	0.45
1:CA:389:A:H2'	1:CA:390:C:H5'	1.98	0.45
1:CA:390:C:O5'	1:CA:390:C:H6	1.98	0.45
1:CA:373:A:C8	1:CA:482:A:C8	3.05	0.45
1:CA:673:G:C6	1:CA:734:G:C6	3.05	0.45
1:CA:666:G:H1'	1:CA:741:G:H22	1.80	0.45
1:CA:895:G:H2'	1:CA:896:C:C6	2.51	0.45
1:CA:928:G:C2	1:CA:1390:U:O2	2.70	0.45
2:CB:178:ARG:HH11	2:CB:178:ARG:CB	2.27	0.45
4:CD:4:TYR:O	4:CD:5:ILE:CB	2.64	0.45
4:CD:4:TYR:O	4:CD:5:ILE:HB	2.16	0.45
4:CD:92:VAL:O	4:CD:96:LEU:HD22	2.16	0.45
5:CE:127:ASN:O	5:CE:128:PRO:C	2.55	0.45
9:CI:33:PHE:C	9:CI:35:GLU:H	2.19	0.45
9:CI:93:ARG:C	9:CI:95:LYS:H	2.19	0.45
11:CK:60:ALA:O	11:CK:62:GLN:N	2.49	0.45
11:CK:72:ALA:HB1	11:CK:77:MET:HE3	1.99	0.45
12:CL:25:PRO:C	12:CL:27:LEU:H	2.19	0.45
12:CL:28:LYS:C	12:CL:30:ALA:N	2.68	0.45
12:CL:89:ARG:HB2	12:CL:89:ARG:CZ	2.46	0.45
1:CA:1228:C:O3'	13:CM:116:THR:HA	2.17	0.45
13:CM:89:GLY:C	13:CM:90:LEU:O	2.55	0.45
16:CP:32:TYR:CE2	16:CP:35:LYS:HB2	2.51	0.45
16:CP:53:VAL:HG23	16:CP:54:GLU:N	2.31	0.45
17:CQ:59:ILE:CD1	17:CQ:73:VAL:HA	2.47	0.45
20:CT:24:LEU:O	20:CT:25:ARG:C	2.54	0.45
25:CY:18:LEU:C	25:CY:20:VAL:N	2.67	0.45
25:CY:38:LEU:O	25:CY:40:HIS:N	2.44	0.45
27:D1:41:ARG:HH11	27:D1:41:ARG:CG	2.21	0.45
28:D2:22:GLU:O	28:D2:24:LEU:N	2.50	0.45
28:D2:53:LEU:O	28:D2:53:LEU:HG	2.16	0.45
32:D6:37:ARG:O	32:D6:48:VAL:O	2.33	0.45
34:D8:29:LYS:NZ	34:D8:44:LYS:CB	2.79	0.45
35:DA:106:C:H1'	55:DY:2:ARG:CZ	2.41	0.45
35:DA:1131:G:C2	35:DA:1132:A:N7	2.84	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1145:C:H2'	35:DA:1146:C:C6	2.51	0.45
35:DA:2330:G:N2	35:DA:2386:C:C2	2.84	0.45
35:DA:2482:G:H1	47:DQ:53:ALA:HB2	1.82	0.45
35:DA:2513:G:H2'	35:DA:2514:U:C6	2.52	0.45
35:DA:1953:A:H2	35:DA:2549:G:N3	2.14	0.45
35:DA:1786:A:C2	35:DA:2606:C:H1'	2.51	0.45
35:DA:2637:U:H1'	35:DA:2782:G:H22	1.82	0.45
35:DA:2764:A:H2'	35:DA:2766:G:C8	2.52	0.45
35:DA:2840:C:O2'	35:DA:2841:C:H5'	2.17	0.45
35:DA:645:C:O2	35:DA:645:C:C2'	2.64	0.45
36:DB:13:A:N1	36:DB:69:G:O2'	2.50	0.45
37:DC:75:LEU:C	37:DC:75:LEU:HD23	2.37	0.45
38:DD:213:ARG:O	38:DD:214:TRP:C	2.54	0.45
38:DD:69:ARG:C	38:DD:71:ASP:H	2.20	0.45
35:DA:574:C:N3	39:DE:145:LYS:HE2	2.31	0.45
40:DF:45:ARG:CG	40:DF:46:ARG:N	2.79	0.45
40:DF:45:ARG:CD	40:DF:46:ARG:H	2.30	0.45
41:DG:7:LEU:O	41:DG:10:LYS:HB2	2.15	0.45
41:DG:180:PHE:HB2	41:DG:182:LYS:HE3	1.99	0.45
41:DG:40:ASN:OD1	41:DG:41:GLN:N	2.50	0.45
42:DH:158:HIS:HE1	42:DH:168:PRO:HB2	1.82	0.45
43:DI:81:VAL:CG2	43:DI:142:VAL:HG13	2.41	0.45
46:DP:113:LYS:HG2	46:DP:115:LEU:CD2	2.46	0.45
46:DP:114:ILE:HD11	46:DP:130:PHE:CZ	2.52	0.45
49:DS:28:VAL:CG1	49:DS:29:PHE:H	2.09	0.45
50:DT:28:VAL:HG21	50:DT:46:GLU:OE2	2.17	0.45
50:DT:52:ILE:HG22	50:DT:61:PHE:HB3	1.97	0.45
51:DU:92:ARG:CG	51:DU:95:LEU:H	2.18	0.45
52:DV:71:LEU:CD1	52:DV:72:VAL:N	2.80	0.45
53:DW:12:ILE:HG13	53:DW:42:ARG:NH1	2.31	0.45
53:DW:45:TYR:CD2	53:DW:45:TYR:C	2.89	0.45
54:DX:35:THR:O	54:DX:36:LYS:O	2.34	0.45
35:DA:143:G:O4'	54:DX:38:GLU:HG3	2.17	0.45
54:DX:62:LYS:HB3	54:DX:69:TYR:H	1.82	0.45
55:DY:88:LYS:CD	55:DY:88:LYS:N	2.78	0.45
1:AA:1303:C:C2'	1:AA:1303:C:O2	2.59	0.45
1:AA:1421:G:H2'	1:AA:1422:G:C8	2.51	0.45
1:AA:136:C:H42	1:AA:227:G:H1	1.65	0.45
1:AA:299:G:C6	1:AA:300:A:N1	2.85	0.45
1:AA:363:A:O2'	1:AA:364:A:H5'	2.16	0.45
1:AA:460:G:N2	1:AA:472:A:H62	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:691:G:O2'	1:AA:692:U:H5'	2.17	0.45
4:AD:172:PRO:HD2	4:AD:173:TRP:CZ3	2.52	0.45
4:AD:187:ARG:NH1	4:AD:187:ARG:HG2	2.29	0.45
7:AG:22:LEU:HD12	7:AG:22:LEU:O	2.16	0.45
9:AI:33:PHE:C	9:AI:35:GLU:H	2.19	0.45
9:AI:47:LEU:N	9:AI:47:LEU:CD1	2.80	0.45
14:AN:8:GLU:O	14:AN:11:LYS:N	2.44	0.45
18:AR:31:LEU:HG	18:AR:65:ILE:HD13	1.99	0.45
18:AR:73:ALA:HB3	18:AR:79:LEU:HD12	1.97	0.45
23:AW:31:G:C5	23:AW:32:G:N7	2.85	0.45
1:AA:1397:C:N4	24:AX:22:U:C4	2.81	0.45
25:AY:108:GLU:O	25:AY:111:ARG:N	2.43	0.45
26:B0:37:LEU:O	26:B0:38:VAL:HG22	2.17	0.45
27:B1:32:LYS:HG2	35:BA:2396:G:O2'	2.17	0.45
27:B1:71:TYR:O	27:B1:75:GLU:HG2	2.16	0.45
28:B2:13:ALA:N	28:B2:14:ARG:HE	2.15	0.45
28:B2:17:SER:O	28:B2:18:PRO:C	2.53	0.45
34:B8:56:GLU:O	34:B8:58:ILE:N	2.50	0.45
34:B8:6:THR:O	34:B8:8:LYS:N	2.50	0.45
35:BA:1303:G:N2	35:BA:1304:C:C2	2.84	0.45
35:BA:1528(A):A:N7	35:BA:1529:G:C8	2.85	0.45
35:BA:1310:G:H1	35:BA:1604:C:H42	1.65	0.45
35:BA:1745(A):C:H5'	35:BA:1746:G:OP2	2.16	0.45
35:BA:2009:G:C2	35:BA:2010:G:C8	3.05	0.45
35:BA:2308:G:N7	35:BA:2310:A:H5'	2.31	0.45
35:BA:2298:A:C2	35:BA:2321:G:C4	3.05	0.45
35:BA:2474:C:H2'	35:BA:2474:C:O2	2.16	0.45
35:BA:2759:G:H5'	35:BA:2759:G:H8	1.81	0.45
35:BA:2762:G:C3'	35:BA:2763:G:C5'	2.94	0.45
35:BA:2792:G:N3	35:BA:2792:G:H2'	2.31	0.45
35:BA:296:C:N4	35:BA:343:C:H42	2.15	0.45
35:BA:610:G:C6	35:BA:611:C:N4	2.85	0.45
35:BA:791:C:O2	35:BA:794:G:H5'	2.16	0.45
35:BA:804:A:H2'	35:BA:806:C:N4	2.31	0.45
36:BB:28:C:H42	36:BB:56:G:H1	1.62	0.45
38:BD:247:ALA:HB2	38:BD:253:GLN:HA	1.99	0.45
35:BA:1655:A:H4'	39:BE:115:GLY:N	2.30	0.45
40:BF:20:LEU:HD12	40:BF:199:TRP:HH2	1.82	0.45
40:BF:20:LEU:HD13	40:BF:203:GLN:OE1	2.17	0.45
41:BG:41:GLN:O	41:BG:43:LEU:N	2.45	0.45
43:BI:98:ALA:HB1	43:BI:109:ILE:CG1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BI:39:ALA:O	43:BI:40:THR:C	2.55	0.45
44:BN:26:LEU:CD2	44:BN:30:ILE:HD11	2.47	0.45
44:BN:33:LEU:HD23	44:BN:38:HIS:CE1	2.52	0.45
46:BP:77:ARG:CB	46:BP:78:PRO:HD2	2.38	0.45
46:BP:96:THR:HG22	46:BP:126:VAL:H	1.82	0.45
47:BQ:43:THR:HG1	47:BQ:45:GLN:HB2	1.80	0.45
47:BQ:75:THR:HA	47:BQ:88:GLY:CA	2.28	0.45
49:BS:101:LEU:HD13	49:BS:101:LEU:O	2.16	0.45
50:BT:115:ARG:HB3	50:BT:116:ALA:H	1.65	0.45
50:BT:33:LYS:HE2	50:BT:43:GLN:OE1	2.17	0.45
50:BT:58:ASN:ND2	50:BT:58:ASN:O	2.50	0.45
55:BY:19:LYS:HG2	55:BY:19:LYS:O	2.17	0.45
55:BY:49:VAL:O	55:BY:53:PRO:CG	2.65	0.45
56:BZ:71:VAL:HG13	56:BZ:86:VAL:CG1	2.47	0.45
56:BZ:28:MET:HA	56:BZ:88:PHE:HB2	1.99	0.45
1:CA:1235:U:O3'	21:CU:3:LYS:HB2	2.17	0.45
1:CA:1380:U:O2	7:CG:3:ARG:NH1	2.46	0.45
1:CA:1387:G:C6	1:CA:1388:C:N4	2.85	0.45
1:CA:259:G:H2'	1:CA:260:G:H8	1.80	0.45
1:CA:302:G:H2'	1:CA:303:A:O4'	2.16	0.45
1:CA:328:C:O2'	1:CA:329:A:OP2	2.28	0.45
1:CA:512:U:H2'	1:CA:513:C:H6	1.81	0.45
1:CA:936:C:H2'	1:CA:937:A:O4'	2.16	0.45
1:CA:939:G:C4	1:CA:940:C:C5	3.04	0.45
2:CB:47:THR:O	2:CB:51:LEU:HG	2.17	0.45
5:CE:89:ILE:CD1	5:CE:91:LEU:HD11	2.46	0.45
7:CG:66:VAL:C	7:CG:68:ASN:H	2.19	0.45
9:CI:118:LYS:HB2	9:CI:121:ARG:HB2	1.99	0.45
12:CL:6:THR:HG23	12:CL:9:GLN:N	2.11	0.45
16:CP:19:ILE:CG2	16:CP:36:ILE:HG13	2.42	0.45
19:CS:10:PHE:CD1	19:CS:10:PHE:N	2.84	0.45
1:CA:263:A:OP2	20:CT:79:ARG:NH1	2.50	0.45
25:CY:121:TYR:O	25:CY:124:GLU:HB2	2.17	0.45
27:D1:23:LYS:NZ	27:D1:23:LYS:CA	2.79	0.45
27:D1:89:GLU:C	27:D1:93:GLU:OE2	2.54	0.45
31:D5:40:LYS:NZ	31:D5:50:GLY:HA2	2.30	0.45
34:D8:40:GLU:CD	34:D8:44:LYS:HE3	2.37	0.45
35:DA:1051:G:C2'	35:DA:1052:C:H5''	2.47	0.45
35:DA:1165:U:O2'	35:DA:1166:C:H5'	2.16	0.45
35:DA:1517:G:C2'	35:DA:1518:U:H5'	2.47	0.45
35:DA:1609:A:C5	35:DA:1616:A:C8	3.05	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1778:U:C5	35:DA:1784:A:N3	2.84	0.45
35:DA:1862:G:C2	35:DA:1863:G:C8	3.05	0.45
35:DA:2057:A:O2'	35:DA:2058:A:H5'	2.17	0.45
35:DA:2415:G:O3'	46:DP:66:GLY:CA	2.50	0.45
35:DA:361:G:C3'	35:DA:362:U:H5''	2.47	0.45
35:DA:426:C:HO2'	35:DA:427:U:H5'	1.82	0.45
35:DA:542:C:C4	35:DA:543:C:N4	2.82	0.45
35:DA:987:G:H2'	35:DA:988:A:O4'	2.17	0.45
38:DD:35:LYS:HZ1	38:DD:104:TYR:HB2	1.80	0.45
35:DA:2579:C:H4'	39:DE:134:ILE:CD1	2.47	0.45
43:DI:114:LEU:O	43:DI:129:THR:O	2.35	0.45
44:DN:28:THR:O	44:DN:31:ALA:HB3	2.17	0.45
47:DQ:110:THR:HG1	47:DQ:113:GLN:HG3	1.80	0.45
47:DQ:58:PHE:O	47:DQ:59:ARG:C	2.54	0.45
47:DQ:90:VAL:O	47:DQ:90:VAL:HG12	2.16	0.45
49:DS:28:VAL:C	49:DS:89:ARG:HG2	2.37	0.45
50:DT:58:ASN:ND2	50:DT:58:ASN:O	2.50	0.45
50:DT:61:PHE:CZ	50:DT:76:PHE:HB3	2.52	0.45
45:DO:76:ALA:HB3	50:DT:75:ILE:HD13	1.97	0.45
51:DU:76:TYR:O	51:DU:80:ILE:HG12	2.16	0.45
51:DU:96:ALA:C	51:DU:98:LEU:N	2.69	0.45
53:DW:9:TYR:CD2	53:DW:9:TYR:N	2.84	0.45
54:DX:14:SER:O	54:DX:15:GLU:C	2.55	0.45
54:DX:31:HIS:HD2	54:DX:33:LYS:O	2.00	0.45
54:DX:76:ARG:O	54:DX:77:LYS:HB2	2.17	0.45
55:DY:49:VAL:O	55:DY:53:PRO:CG	2.65	0.45
55:DY:65:ALA:O	55:DY:67:LEU:N	2.50	0.45
56:DZ:10:ARG:HG2	56:DZ:38:TYR:HD2	1.81	0.45
56:DZ:89:PHE:CE1	56:DZ:96:VAL:HG21	2.52	0.45
1:AA:1006:C:H2'	1:AA:1007:C:C6	2.52	0.45
1:AA:1105:A:H2'	1:AA:1106:G:H8	1.82	0.45
1:AA:1315:U:O2'	1:AA:1316:G:H5'	2.16	0.45
1:AA:1346:A:H5''	9:AI:120:ARG:NH1	2.24	0.45
1:AA:1423:G:H2'	1:AA:1424:C:C6	2.52	0.45
1:AA:1436:U:O2'	1:AA:1437:C:H5'	2.17	0.45
1:AA:21:G:H2'	1:AA:22:G:C8	2.51	0.45
1:AA:408:A:H2'	1:AA:409:G:O4'	2.15	0.45
1:AA:761:G:H2'	1:AA:762:C:H6	1.81	0.45
1:AA:902:G:O2'	1:AA:903:G:H5'	2.17	0.45
1:AA:936:C:H2'	1:AA:937:A:O4'	2.17	0.45
2:AB:72:GLY:HA2	2:AB:165:VAL:HG22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:88:ALA:CB	2:AB:223:ILE:HD11	2.47	0.45
4:AD:170:VAL:CG1	4:AD:174:LEU:HB2	2.46	0.45
7:AG:102:ARG:HG3	7:AG:106:GLN:NE2	2.30	0.45
7:AG:47:CYS:CA	7:AG:58:PRO:HG3	2.47	0.45
7:AG:69:VAL:HG12	7:AG:69:VAL:O	2.17	0.45
8:AH:120:THR:O	8:AH:122:ARG:N	2.50	0.45
8:AH:45:ILE:HD13	8:AH:62:TYR:O	2.16	0.45
10:AJ:47:PHE:HD1	10:AJ:47:PHE:O	2.00	0.45
11:AK:96:ARG:O	11:AK:99:GLN:HG2	2.16	0.45
13:AM:36:LYS:HB2	13:AM:59:TYR:CZ	2.51	0.45
13:AM:68:GLY:O	13:AM:70:LEU:N	2.49	0.45
15:AO:82:ILE:HG23	15:AO:83:GLU:N	2.32	0.45
16:AP:29:ASP:OD2	16:AP:29:ASP:N	2.50	0.45
1:AA:277:C:OP1	17:AQ:41:LYS:HE3	2.17	0.45
18:AR:43:PHE:CD1	18:AR:43:PHE:N	2.81	0.45
23:AW:40:C:O2'	23:AW:41:C:H5'	2.17	0.45
24:AX:13:A:HO2'	24:AX:14:U:H5'	1.81	0.45
25:AY:6:LEU:C	25:AY:8:ALA:N	2.70	0.45
25:AY:6:LEU:C	25:AY:6:LEU:HD23	2.37	0.45
26:B0:36:ILE:HA	26:B0:60:PHE:CB	2.47	0.45
33:B7:39:ARG:HH11	33:B7:39:ARG:HG2	1.80	0.45
34:B8:22:VAL:HB	34:B8:53:PRO:HB3	1.95	0.45
35:BA:1187:G:OP1	52:BV:82:ARG:NH2	2.42	0.45
35:BA:1213:A:H2'	35:BA:1214:A:C8	2.47	0.45
35:BA:1635:G:H2'	35:BA:1636:C:H6	1.82	0.45
35:BA:1655:A:O2'	39:BE:115:GLY:HA3	2.17	0.45
35:BA:1695:G:H2'	35:BA:1696:G:C4'	2.46	0.45
35:BA:1799:G:O3'	38:BD:264:LYS:NZ	2.47	0.45
35:BA:1910:G:N1	35:BA:1921:G:C5	2.85	0.45
35:BA:1786:A:C2	35:BA:2606:C:H1'	2.51	0.45
35:BA:2820:A:H8	39:BE:191:PRO:CB	2.30	0.45
35:BA:373:U:H1'	35:BA:423:A:N3	2.32	0.45
35:BA:448:U:H4'	35:BA:449:A:OP2	2.17	0.45
35:BA:621:A:H2'	35:BA:622:G:C5'	2.42	0.45
35:BA:832:G:OP1	46:BP:40:SER:HB3	2.16	0.45
36:BB:82:G:H2'	36:BB:83:G:C8	2.49	0.45
38:BD:106:ILE:HD13	38:BD:157:ARG:HB2	1.99	0.45
38:BD:138:VAL:HG21	38:BD:166:GLN:O	2.17	0.45
38:BD:218:ARG:HG3	38:BD:218:ARG:HH11	1.82	0.45
39:BE:163:GLU:O	39:BE:165:VAL:N	2.50	0.45
40:BF:101:LEU:HD12	40:BF:102:PRO:CD	2.39	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:443:A:OP1	40:BF:46:ARG:HB2	2.16	0.45
42:BH:59:ARG:HG2	42:BH:59:ARG:O	2.17	0.45
42:BH:67:LEU:C	42:BH:71:LEU:HD22	2.37	0.45
44:BN:67:LEU:C	44:BN:69:GLN:N	2.69	0.45
46:BP:100:LEU:CD2	46:BP:100:LEU:N	2.79	0.45
46:BP:59:LEU:HA	46:BP:61:ARG:HD2	1.99	0.45
35:BA:631:A:H4'	46:BP:65:ARG:HG3	1.98	0.45
46:BP:80:TYR:CZ	46:BP:111:ARG:HG2	2.52	0.45
47:BQ:66:ILE:HG22	47:BQ:104:PHE:HE2	1.74	0.45
47:BQ:54:MET:HG3	47:BQ:64:ILE:HD13	1.99	0.45
47:BQ:7:MET:O	47:BQ:8:LYS:HB3	2.17	0.45
49:BS:38:GLN:HG3	49:BS:49:VAL:O	2.16	0.45
49:BS:87:PHE:HZ	49:BS:97:ARG:HH21	1.63	0.45
50:BT:106:SER:C	50:BT:107:ASP:OD1	2.55	0.45
50:BT:109:GLU:O	50:BT:110:ILE:C	2.54	0.45
50:BT:29:ARG:HG3	50:BT:84:GLN:O	2.15	0.45
55:BY:15:VAL:HG12	55:BY:17:SER:H	1.81	0.45
56:BZ:108:PRO:C	56:BZ:109:ALA:O	2.52	0.45
56:BZ:28:MET:HG2	56:BZ:37:VAL:HG22	1.99	0.45
1:CA:1030(C):G:N7	1:CA:1031:G:N2	2.65	0.45
1:CA:1465:C:H2'	1:CA:1466:C:H6	1.81	0.45
1:CA:152:A:N6	1:CA:170:U:N3	2.65	0.45
1:CA:18:C:H5''	5:CE:127:ASN:ND2	2.32	0.45
1:CA:190:U:O2'	1:CA:191:G:H5'	2.17	0.45
1:CA:243:A:C2	1:CA:246:A:C8	3.05	0.45
1:CA:245:C:C2'	1:CA:246:A:H5'	2.47	0.45
1:CA:656:C:O2'	1:CA:657:G:H5'	2.17	0.45
1:CA:814:A:C8	1:CA:816:A:C8	3.05	0.45
2:CB:102:LEU:HD22	2:CB:176:GLU:HB3	1.98	0.45
2:CB:23:ARG:HG2	2:CB:23:ARG:O	2.16	0.45
2:CB:22:LYS:O	2:CB:24:TRP:N	2.50	0.45
4:CD:104:VAL:HG21	4:CD:140:VAL:CG2	2.47	0.45
4:CD:173:TRP:C	4:CD:186:LEU:HB2	2.37	0.45
4:CD:178:VAL:C	4:CD:180:GLY:H	2.19	0.45
6:CF:10:LEU:HD12	6:CF:10:LEU:N	2.31	0.45
7:CG:121:ALA:N	7:CG:124:LEU:HD12	2.31	0.45
7:CG:94:ARG:HG3	7:CG:94:ARG:H	1.63	0.45
9:CI:17:VAL:HA	9:CI:63:ILE:HG23	1.98	0.45
11:CK:116:HIS:O	11:CK:117:ASN:HB3	2.17	0.45
11:CK:48:ILE:C	11:CK:50:TYR:H	2.20	0.45
13:CM:71:ARG:O	13:CM:74:VAL:HB	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:973:G:O3'	14:CN:41:ARG:NH2	2.49	0.45
18:CR:44:LEU:O	18:CR:51:LEU:HD21	2.17	0.45
19:CS:22:LEU:C	19:CS:22:LEU:HD13	2.37	0.45
23:CW:51:U:H3	23:CW:65:G:H22	1.64	0.45
26:D0:36:ILE:HA	26:D0:60:PHE:CB	2.47	0.45
28:D2:30:ARG:N	28:D2:30:ARG:CD	2.78	0.45
29:D3:29:ARG:O	29:D3:30:ARG:C	2.55	0.45
32:D6:51:GLU:O	32:D6:52:VAL:HB	2.17	0.45
33:D7:29:LYS:HZ2	33:D7:32:LYS:NZ	2.14	0.45
35:DA:1204:A:N1	35:DA:1241:A:C2	2.85	0.45
35:DA:1753:G:C8	35:DA:1755:A:OP2	2.70	0.45
35:DA:1759:A:H2'	35:DA:1760:A:H8	1.80	0.45
35:DA:2553:G:H3'	35:DA:2554:U:H5''	1.98	0.45
35:DA:2640:G:H5'	35:DA:2640:G:C8	2.51	0.45
35:DA:2725:A:N7	35:DA:2727:G:C5	2.85	0.45
35:DA:2755:C:H6	35:DA:2755:C:H3'	1.81	0.45
35:DA:295:G:C2	35:DA:296:C:C6	3.05	0.45
35:DA:29:U:O2'	35:DA:30:G:H5'	2.16	0.45
35:DA:307:G:N2	35:DA:309:G:H3'	2.32	0.45
35:DA:374:A:H2'	35:DA:375:C:O4'	2.17	0.45
35:DA:479:A:HO2'	35:DA:481:G:H8	1.63	0.45
35:DA:718:A:H2'	35:DA:719:C:O4'	2.16	0.45
35:DA:869:G:O2'	35:DA:870:A:H5'	2.16	0.45
36:DB:74:U:C3'	36:DB:75:G:C5'	2.90	0.45
38:DD:21:PHE:O	38:DD:22:SER:C	2.54	0.45
39:DE:173:VAL:CG1	39:DE:174:ASP:N	2.80	0.45
39:DE:48:GLN:CG	39:DE:78:LEU:HD12	2.45	0.45
40:DF:1:MET:O	40:DF:3:GLU:N	2.49	0.45
35:DA:443:A:OP1	40:DF:46:ARG:HB2	2.17	0.45
41:DG:107:LEU:HD13	41:DG:178:PHE:HE1	1.82	0.45
41:DG:104:GLU:O	41:DG:108:ASN:N	2.50	0.45
42:DH:66:GLY:O	42:DH:67:LEU:C	2.54	0.45
44:DN:17:ASP:OD2	44:DN:55:VAL:O	2.34	0.45
44:DN:65:LYS:C	44:DN:66:LYS:HE3	2.37	0.45
47:DQ:85:LYS:HG3	47:DQ:86:GLY:H	1.79	0.45
48:DR:76:VAL:HG13	48:DR:77:ARG:N	2.31	0.45
50:DT:80:SER:CB	50:DT:81:PRO:HD2	2.35	0.45
51:DU:26:GLY:C	51:DU:28:ARG:H	2.20	0.45
51:DU:52:ARG:O	51:DU:55:ARG:N	2.49	0.45
53:DW:50:VAL:HG13	53:DW:105:VAL:HG21	1.99	0.45
54:DX:89:ILE:O	54:DX:89:ILE:CG2	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:61:ILE:O	55:DY:62:GLU:HB2	2.16	0.45
56:DZ:175:VAL:HA	56:DZ:176:PRO:HD3	1.84	0.45
1:AA:1197:G:O2'	1:AA:1198:G:H5'	2.17	0.45
1:AA:1316:G:H2'	1:AA:1317:C:H5'	1.99	0.45
1:AA:276:G:O2'	1:AA:277:C:H5'	2.17	0.45
1:AA:327:A:C2	1:AA:329:A:C4	3.05	0.45
1:AA:359:U:H2'	1:AA:360:A:H8	1.80	0.45
1:AA:63:C:H42	1:AA:104:G:H1	1.64	0.45
1:AA:863:U:H2'	1:AA:865:A:OP2	2.16	0.45
1:AA:922:G:H4'	5:AE:20:GLN:HA	1.99	0.45
2:AB:236:TYR:HB3	2:AB:239:VAL:HB	1.99	0.45
2:AB:47:THR:O	2:AB:51:LEU:HG	2.16	0.45
2:AB:74:LYS:O	2:AB:75:LYS:C	2.55	0.45
3:AC:112:SER:OG	3:AC:114:PRO:HD2	2.17	0.45
3:AC:76:VAL:CG2	3:AC:77:ILE:N	2.79	0.45
4:AD:21:LEU:HD22	4:AD:115:ARG:HG3	1.98	0.45
1:AA:18:C:C5'	5:AE:127:ASN:HD21	2.28	0.45
5:AE:147:ASP:C	5:AE:150:ARG:HB3	2.37	0.45
5:AE:77:PRO:HG2	5:AE:78:HIS:H	1.82	0.45
6:AF:45:LEU:CD1	6:AF:46:ARG:N	2.78	0.45
6:AF:98:LEU:HD12	6:AF:98:LEU:N	2.31	0.45
7:AG:120:ILE:HG22	7:AG:124:LEU:CD1	2.46	0.45
7:AG:50:ILE:O	7:AG:50:ILE:HG22	2.16	0.45
7:AG:57:GLU:C	7:AG:59:LEU:H	2.20	0.45
8:AH:26:VAL:CG1	8:AH:59:LEU:HB2	2.47	0.45
9:AI:119:ALA:O	9:AI:120:ARG:HB2	2.17	0.45
9:AI:89:ASN:C	9:AI:91:ASP:H	2.19	0.45
11:AK:18:ARG:O	11:AK:33:THR:N	2.47	0.45
11:AK:41:THR:HG21	11:AK:71:LYS:CD	2.47	0.45
15:AO:32:LEU:O	15:AO:36:ILE:HG13	2.16	0.45
15:AO:53:HIS:O	15:AO:56:LEU:HB3	2.16	0.45
17:AQ:59:ILE:HG23	17:AQ:71:PHE:HB3	1.98	0.45
18:AR:40:LEU:O	18:AR:42:ARG:N	2.50	0.45
25:AY:172:ALA:O	25:AY:175:LEU:N	2.49	0.45
33:B7:48:LYS:CD	33:B7:48:LYS:N	2.79	0.45
34:B8:35:GLN:NE2	34:B8:36:LYS:HZ2	2.15	0.45
35:BA:1224:C:O2'	35:BA:1225:G:H5'	2.16	0.45
35:BA:1557:C:OP2	35:BA:1558:A:O2'	2.35	0.45
35:BA:1578:U:H2'	35:BA:1579:A:H5''	1.99	0.45
35:BA:184:C:H2'	35:BA:185:U:H6	1.77	0.45
35:BA:1885:A:H2'	35:BA:1886:C:O4'	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2014:A:C2	35:BA:2015:A:C2	3.05	0.45
35:BA:2037:G:H2'	35:BA:2038:G:H8	1.79	0.45
35:BA:2041:U:H2'	35:BA:2042:A:C8	2.52	0.45
35:BA:2050:C:H2'	35:BA:2051:A:C8	2.51	0.45
35:BA:2070:G:C2	35:BA:2071:A:C4	3.05	0.45
26:B0:16:SER:CB	35:BA:2261:C:H3'	2.47	0.45
35:BA:2552:U:H2'	35:BA:2554:U:H5''	1.98	0.45
35:BA:2683:C:O2'	35:BA:2684:U:H5'	2.16	0.45
35:BA:2689:U:O2'	35:BA:2690:C:OP2	2.26	0.45
35:BA:2738:A:C2	35:BA:2739:U:C6	3.05	0.45
35:BA:361:G:C3'	35:BA:362:U:H5''	2.46	0.45
35:BA:869:G:O2'	47:BQ:8:LYS:HD3	2.17	0.45
36:BB:60:C:O2'	36:BB:61:G:H5'	2.16	0.45
38:BD:133:LEU:O	38:BD:136:ILE:HG13	2.16	0.45
38:BD:198:ASN:ND2	38:BD:198:ASN:C	2.69	0.45
40:BF:103:LYS:HA	40:BF:106:ARG:HE	1.81	0.45
41:BG:161:THR:HG22	41:BG:163:ALA:N	2.31	0.45
44:BN:103:VAL:O	44:BN:106:MET:HB2	2.16	0.45
45:BO:6:THR:CG2	45:BO:7:TYR:H	2.23	0.45
47:BQ:20:ALA:O	47:BQ:23:GLY:N	2.37	0.45
50:BT:66:VAL:HG13	50:BT:71:GLY:N	2.31	0.45
51:BU:105:VAL:HG12	51:BU:109:LEU:HD11	1.99	0.45
35:BA:996:A:H4'	51:BU:92:ARG:NE	2.32	0.45
55:BY:71:LYS:HZ2	55:BY:71:LYS:HB2	1.82	0.45
1:CA:10:A:H2'	1:CA:11:G:C8	2.51	0.45
1:CA:1105:A:H2'	1:CA:1106:G:H8	1.81	0.45
1:CA:1231:G:O2'	1:CA:1232:U:H5'	2.16	0.45
1:CA:1277:C:H2'	1:CA:1278:U:H5'	1.99	0.45
1:CA:1351:U:H6	1:CA:1351:U:O5'	2.00	0.45
1:CA:1425:U:H1'	1:CA:1476:G:N2	2.32	0.45
1:CA:502:G:C2	1:CA:544:G:C2	3.05	0.45
1:CA:514:C:O2'	1:CA:515:G:H5'	2.17	0.45
1:CA:780:A:C2	1:CA:803:G:C6	3.05	0.45
1:CA:920:U:C2	1:CA:921:U:C5	3.05	0.45
2:CB:236:TYR:HB3	2:CB:239:VAL:HB	1.99	0.45
2:CB:69:LEU:C	2:CB:69:LEU:CD1	2.85	0.45
3:CC:178:LEU:C	3:CC:180:ALA:H	2.19	0.45
1:CA:932:C:H4'	7:CG:4:ARG:NH2	2.32	0.45
8:CH:107:LEU:HD23	8:CH:107:LEU:O	2.17	0.45
8:CH:10:LEU:O	8:CH:14:ARG:HB2	2.17	0.45
8:CH:5:PRO:O	8:CH:8:ASP:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CK:64:ALA:C	11:CK:66:LEU:N	2.69	0.45
11:CK:41:THR:HG21	11:CK:71:LYS:CD	2.47	0.45
12:CL:32:PHE:CB	12:CL:84:LEU:HD21	2.46	0.45
19:CS:10:PHE:HE2	19:CS:70:LYS:NZ	2.15	0.45
21:CU:24:ARG:HD2	21:CU:24:ARG:N	2.32	0.45
23:CW:43:G:H2'	23:CW:44:A:C5'	2.46	0.45
25:CY:64:ARG:HD2	25:CY:64:ARG:N	2.31	0.45
26:D0:7:LEU:CA	47:DQ:83:MET:SD	3.05	0.45
33:D7:2:LYS:HG2	35:DA:1620:G:O2'	2.17	0.45
35:DA:1013:C:O2'	35:DA:1014:U:H5'	2.16	0.45
35:DA:1350:C:C2'	35:DA:1351:C:H5'	2.47	0.45
35:DA:1531:C:H3'	35:DA:1532:C:C5'	2.46	0.45
35:DA:1638:C:H1'	35:DA:2698:U:O2'	2.17	0.45
35:DA:202:U:O2'	35:DA:203:C:H5'	2.16	0.45
35:DA:2228:G:C6	35:DA:2229:C:C4	3.05	0.45
35:DA:2376:A:H2'	35:DA:2377:A:O4'	2.17	0.45
35:DA:2711:A:OP2	35:DA:2712(A):A:OP2	2.35	0.45
35:DA:2740:A:C2'	35:DA:2741:A:C8	2.92	0.45
35:DA:27:G:H2'	35:DA:28:A:OP2	2.17	0.45
35:DA:2861:G:C2	35:DA:2862:G:C8	3.05	0.45
35:DA:469:G:C2'	35:DA:470:A:H5''	2.46	0.45
35:DA:523:C:C5	35:DA:524:U:H5	2.34	0.45
35:DA:619:G:H3'	35:DA:620:G:N2	2.31	0.45
35:DA:674:G:H2'	35:DA:804:A:H61	1.82	0.45
35:DA:684:G:N2	35:DA:787:U:H2'	2.31	0.45
35:DA:808:G:O2'	35:DA:1254:A:O2'	2.30	0.45
35:DA:904:C:H2'	35:DA:905:U:H5'	1.99	0.45
35:DA:949:C:O2'	35:DA:950:G:H5'	2.16	0.45
36:DB:7:G:H2'	36:DB:8:U:O4'	2.16	0.45
38:DD:247:ALA:CB	38:DD:253:GLN:HA	2.46	0.45
38:DD:85:ASP:OD1	38:DD:87:ASN:HB2	2.17	0.45
39:DE:119:ARG:HD2	39:DE:120:TRP:CE2	2.51	0.45
39:DE:83:ASP:O	39:DE:84:PHE:HB2	2.17	0.45
35:DA:443:A:OP1	40:DF:46:ARG:CB	2.64	0.45
41:DG:39:ILE:HA	41:DG:157:ILE:CB	2.46	0.45
41:DG:180:PHE:CB	41:DG:182:LYS:HG3	2.47	0.45
42:DH:136:ILE:H	42:DH:136:ILE:CD1	2.23	0.45
42:DH:151:ILE:N	42:DH:151:ILE:HD13	2.32	0.45
43:DI:12:LEU:N	43:DI:12:LEU:HD23	2.32	0.45
45:DO:44:LYS:HD3	45:DO:44:LYS:HA	1.81	0.45
46:DP:23:PRO:HB2	46:DP:33:ARG:NE	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:95:VAL:HG23	46:DP:125:VAL:CB	2.46	0.45
35:DA:2467:C:H4'	47:DQ:123:HIS:CE1	2.52	0.45
49:DS:89:ARG:CB	49:DS:97:ARG:HH22	2.30	0.45
50:DT:62:THR:CG2	50:DT:75:ILE:HA	2.24	0.45
51:DU:87:GLY:O	51:DU:88:ILE:HG23	2.16	0.45
53:DW:86:LEU:HA	53:DW:87:PRO:HD3	1.81	0.45
54:DX:82:GLN:HG3	54:DX:83:VAL:H	1.74	0.45
56:DZ:100:VAL:HG22	56:DZ:101:PRO:HD2	1.99	0.45
1:AA:1040:U:H2'	1:AA:1041:A:H8	1.81	0.45
1:AA:1046:A:H2'	1:AA:1046:A:N3	2.30	0.45
1:AA:1106:G:H5''	3:AC:172:ARG:HG2	1.98	0.45
1:AA:1302:U:O2'	1:AA:1303:C:H5'	2.17	0.45
1:AA:1342:C:H2'	1:AA:1343:G:C8	2.51	0.45
1:AA:296:U:H2'	1:AA:297:G:O4'	2.17	0.45
1:AA:451:A:C1'	1:AA:452:A:C8	2.99	0.45
1:AA:685:G:H21	1:AA:686:U:H3	1.61	0.45
1:AA:833:U:H2'	1:AA:834:C:C5	2.52	0.45
1:AA:662:G:O2'	1:AA:836:G:H5''	2.17	0.45
1:AA:972:C:H5'	10:AJ:57:LYS:HZ3	1.82	0.45
2:AB:114:ARG:HD2	2:AB:118:LEU:HG	1.99	0.45
2:AB:197:VAL:HB	2:AB:200:ILE:HG12	1.99	0.45
2:AB:221:LEU:C	2:AB:221:LEU:HD13	2.37	0.45
1:AA:1057:G:H5''	3:AC:154:SER:HG	1.80	0.45
4:AD:108:LEU:N	4:AD:108:LEU:HD12	2.32	0.45
4:AD:128:VAL:CG1	4:AD:129:ASN:H	2.12	0.45
4:AD:61:LYS:NZ	4:AD:62:GLN:NE2	2.65	0.45
6:AF:10:LEU:HD13	6:AF:61:LEU:CD1	2.45	0.45
6:AF:69:GLU:O	6:AF:70:ASP:C	2.55	0.45
9:AI:125:TYR:CD2	9:AI:125:TYR:C	2.88	0.45
11:AK:69:ALA:O	11:AK:73:MET:HG2	2.17	0.45
16:AP:40:ASP:H	16:AP:48:TRP:HB2	1.82	0.45
17:AQ:10:VAL:HG12	17:AQ:53:LEU:CD1	2.46	0.45
18:AR:44:LEU:O	18:AR:51:LEU:HD21	2.17	0.45
28:B2:32:LEU:HG	28:B2:33:MET:H	1.81	0.45
28:B2:46:GLN:O	28:B2:50:ILE:O	2.35	0.45
31:B5:23:HIS:O	31:B5:24:ALA:C	2.56	0.45
35:BA:1175:U:O3'	35:BA:1176:G:H3'	2.16	0.45
35:BA:814:C:H1'	35:BA:1225:G:H21	1.81	0.45
35:BA:1406:U:H3'	35:BA:1407:C:C6	2.46	0.45
35:BA:151:C:H2'	35:BA:152:G:H8	1.80	0.45
35:BA:1444:G:C2	35:BA:1548:C:C2	3.05	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1710:C:H2'	35:BA:1711:C:H6	1.80	0.45
35:BA:178:G:O2'	35:BA:179:G:H5'	2.17	0.45
35:BA:2247:A:H2'	35:BA:2248:C:H6	1.81	0.45
35:BA:2350:C:H2'	35:BA:2351:G:O4'	2.17	0.45
26:B0:39:ARG:NH2	35:BA:2354:G:N2	2.64	0.45
35:BA:239:U:O2'	35:BA:622:G:H4'	2.16	0.45
35:BA:2627:G:N2	35:BA:2781:A:C2	2.85	0.45
35:BA:27:G:C2'	35:BA:28:A:OP2	2.64	0.45
35:BA:2831:G:P	39:BE:58:ARG:NH1	2.90	0.45
35:BA:300:A:C6	35:BA:334:C:H5'	2.52	0.45
35:BA:447:A:C4	35:BA:473:G:C8	3.05	0.45
35:BA:523:C:C4	35:BA:524:U:C5	3.05	0.45
35:BA:576:U:H2'	35:BA:577:G:H8	1.82	0.45
35:BA:702:G:C2	35:BA:731:C:N3	2.85	0.45
35:BA:838:C:N4	35:BA:940:G:H1	2.14	0.45
35:BA:994:C:H1'	52:BV:10:LYS:HE2	1.99	0.45
36:BB:74:U:H3'	36:BB:75:G:H5''	1.97	0.45
38:BD:206:LEU:HD23	38:BD:211:ARG:NH1	2.32	0.45
40:BF:178:PRO:HG2	40:BF:179:GLU:HG3	1.99	0.45
40:BF:34:TRP:O	40:BF:35:GLU:C	2.54	0.45
35:BA:1256:G:O2'	40:BF:82:ILE:HD12	2.16	0.45
42:BH:66:GLY:O	42:BH:69:ARG:N	2.50	0.45
43:BI:12:LEU:HD23	43:BI:12:LEU:N	2.32	0.45
44:BN:76:SER:O	44:BN:77:GLY:O	2.34	0.45
46:BP:114:ILE:HD11	46:BP:130:PHE:CE2	2.52	0.45
46:BP:16:ARG:HD3	46:BP:16:ARG:O	2.17	0.45
47:BQ:110:THR:HG1	47:BQ:113:GLN:HG3	1.81	0.45
49:BS:101:LEU:C	49:BS:101:LEU:HD13	2.38	0.45
49:BS:89:ARG:CB	49:BS:97:ARG:HH22	2.30	0.45
50:BT:83:ILE:CG1	50:BT:84:GLN:HG2	2.45	0.45
52:BV:2:PHE:O	52:BV:3:ALA:HB2	2.16	0.45
52:BV:72:VAL:O	52:BV:73:SER:HB3	2.17	0.45
52:BV:85:LYS:O	52:BV:87:HIS:N	2.45	0.45
35:BA:1601:G:OP2	54:BX:58:HIS:CD2	2.70	0.45
54:BX:63:LYS:O	54:BX:68:ARG:HA	2.17	0.45
56:BZ:125:LEU:O	56:BZ:165:VAL:CG2	2.65	0.45
1:CA:189(E):U:O2'	1:CA:189(F):U:H5'	2.16	0.45
1:CA:192:U:C2	1:CA:193:C:C5	3.05	0.45
1:CA:602:A:O2'	1:CA:603:U:H5'	2.17	0.45
1:CA:802:A:C2'	1:CA:803:G:H5'	2.46	0.45
1:CA:1075:C:H5'	2:CB:103:THR:HG21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:58:LEU:CD1	4:CD:59:ARG:N	2.77	0.45
4:CD:63:LYS:O	4:CD:64:LEU:C	2.54	0.45
5:CE:145:LYS:C	5:CE:148:VAL:HB	2.36	0.45
5:CE:20:GLN:O	5:CE:23:GLY:O	2.35	0.45
5:CE:55:VAL:O	5:CE:58:ALA:HB3	2.16	0.45
6:CF:72:VAL:O	6:CF:75:LEU:HB2	2.17	0.45
8:CH:23:SER:HB2	8:CH:61:VAL:O	2.16	0.45
9:CI:25:LYS:HG3	9:CI:25:LYS:O	2.16	0.45
9:CI:47:LEU:N	9:CI:47:LEU:CD1	2.79	0.45
11:CK:48:ILE:O	11:CK:50:TYR:N	2.42	0.45
12:CL:119:LYS:HD3	12:CL:120:TYR:CE1	2.50	0.45
14:CN:37:PHE:CE1	14:CN:53:LEU:HD22	2.50	0.45
15:CO:56:LEU:O	15:CO:57:LEU:C	2.55	0.45
23:CW:19:G:H1'	23:CW:59:A:C2	2.51	0.45
25:CY:150:SER:C	25:CY:152:ASP:N	2.71	0.45
26:D0:51:VAL:CG2	26:D0:81:VAL:HG23	2.47	0.45
28:D2:36:ARG:NH1	54:DX:92:LEU:HD22	2.32	0.45
29:D3:40:THR:O	29:D3:43:ILE:N	2.49	0.45
34:D8:59:LYS:CB	34:D8:59:LYS:NZ	2.75	0.45
35:DA:121:G:H2'	35:DA:122:G:C8	2.52	0.45
35:DA:154:G:C6	35:DA:173:G:C6	3.05	0.45
35:DA:1935:G:H1'	35:DA:1964:G:H21	1.82	0.45
35:DA:1989:G:H2'	35:DA:1990:C:H5'	1.98	0.45
35:DA:2024:G:H2'	35:DA:2025:C:C6	2.50	0.45
35:DA:2092:U:C4'	35:DA:2093:G:H5''	2.37	0.45
35:DA:2474:C:C2'	35:DA:2474:C:O2	2.64	0.45
35:DA:2746:U:H2'	35:DA:2747:G:H8	1.81	0.45
35:DA:470:A:H2'	35:DA:471:A:O4'	2.17	0.45
35:DA:479:A:H61	35:DA:503:A:H61	1.65	0.45
35:DA:61:G:H1	35:DA:94:C:N4	2.13	0.45
35:DA:702:G:C2	35:DA:731:C:N3	2.85	0.45
35:DA:79:G:C4	35:DA:80:G:C8	3.05	0.45
35:DA:916:G:H2'	35:DA:917:A:H5''	1.99	0.45
36:DB:16:G:H2'	36:DB:17:C:C6	2.51	0.45
37:DC:42:GLU:O	37:DC:212:VAL:HA	2.17	0.45
39:DE:132:HIS:CD2	39:DE:135:HIS:NE2	2.84	0.45
39:DE:72:VAL:O	39:DE:73:GLU:O	2.34	0.45
39:DE:88:GLY:O	39:DE:89:ASP:CB	2.64	0.45
43:DI:89:TYR:O	43:DI:121:LYS:HE2	2.17	0.45
43:DI:87:LYS:HD2	43:DI:121:LYS:O	2.17	0.45
44:DN:29:LYS:O	44:DN:31:ALA:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:77:GLY:O	44:DN:78:TYR:CB	2.64	0.45
45:DO:11:ALA:HB2	45:DO:64:ARG:NH2	2.32	0.45
49:DS:27:SER:CA	49:DS:89:ARG:HD2	2.42	0.45
49:DS:97:ARG:O	49:DS:98:VAL:C	2.54	0.45
53:DW:103:ILE:HG13	53:DW:103:ILE:O	2.17	0.45
53:DW:65:LEU:CD2	53:DW:67:ASP:HB2	2.47	0.45
53:DW:65:LEU:HD22	53:DW:68:ARG:H	1.81	0.45
54:DX:51:VAL:CG1	54:DX:80:ILE:N	2.79	0.45
56:DZ:69:THR:O	56:DZ:70:LEU:HD23	2.17	0.45
1:AA:1203:C:H2'	1:AA:1204:A:O4'	2.18	0.45
1:AA:1293:G:O2'	1:AA:1294:G:P	2.75	0.45
1:AA:1476:G:C2	1:AA:1477:C:C2	3.05	0.45
1:AA:553:A:O2'	1:AA:554:C:H5'	2.16	0.45
1:AA:61:G:O2'	1:AA:62:U:H5'	2.17	0.45
1:AA:690:G:N1	1:AA:691:G:C2	2.85	0.45
1:AA:785:G:C2'	1:AA:786:G:H5'	2.47	0.45
1:AA:586:C:H1'	1:AA:878:G:O2'	2.17	0.45
2:AB:105:PHE:O	2:AB:106:LYS:C	2.55	0.45
2:AB:236:TYR:C	2:AB:238:LEU:H	2.20	0.45
7:AG:6:ARG:O	7:AG:7:ALA:C	2.55	0.45
8:AH:7:ALA:HA	8:AH:10:LEU:HG	1.98	0.45
1:AA:1118:C:P	9:AI:104:ARG:HG3	2.57	0.45
10:AJ:8:LEU:HB2	10:AJ:16:LEU:HD11	1.98	0.45
13:AM:8:GLU:C	13:AM:9:ILE:HD12	2.38	0.45
17:AQ:12:SER:HA	17:AQ:14:LYS:HZ1	1.82	0.45
18:AR:36:ASN:HB2	18:AR:39:VAL:CG2	2.46	0.45
19:AS:41:VAL:CG1	19:AS:44:MET:HB2	2.46	0.45
20:AT:64:ASP:OD1	20:AT:81:LYS:HD2	2.17	0.45
25:AY:142:LYS:HD2	25:AY:142:LYS:O	2.17	0.45
26:B0:12:ASN:O	26:B0:14:ARG:N	2.50	0.45
28:B2:53:LEU:HG	35:BA:77:C:H5'	1.97	0.45
29:B3:22:ALA:O	29:B3:25:ALA:HB3	2.16	0.45
31:B5:40:LYS:HZ3	31:B5:46:CYS:N	2.14	0.45
32:B6:32:ASN:CG	32:B6:33:LYS:N	2.64	0.45
34:B8:25:MET:HB3	34:B8:26:LYS:H	1.52	0.45
34:B8:30:ARG:O	34:B8:30:ARG:CG	2.65	0.45
35:BA:1230:C:O2'	35:BA:1231:G:H5'	2.17	0.45
35:BA:1264:G:C3'	35:BA:1265:A:H5''	2.47	0.45
35:BA:1281:G:O6	35:BA:1286:A:N7	2.50	0.45
35:BA:1301:A:O2'	35:BA:1302:A:P	2.75	0.45
35:BA:1753:G:C8	35:BA:1755:A:OP2	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1798:U:N3	35:BA:1819:A:C2	2.85	0.45
35:BA:16:G:H2'	35:BA:17:G:H8	1.82	0.45
35:BA:1831:G:H2'	35:BA:1832:C:C6	2.52	0.45
27:B1:41:ARG:HH12	35:BA:189:G:C3'	2.30	0.45
35:BA:2350:C:O2'	35:BA:2351:G:H5'	2.16	0.45
35:BA:2474:C:O2	35:BA:2474:C:C2'	2.65	0.45
35:BA:2714:G:H8	35:BA:2714:G:H5''	1.81	0.45
35:BA:27:G:H2'	35:BA:28:A:OP2	2.17	0.45
35:BA:491:G:N2	35:BA:492:A:H1'	2.32	0.45
35:BA:259:G:N2	35:BA:621:A:H8	2.07	0.45
35:BA:967:C:C2'	35:BA:968:G:H5'	2.47	0.45
36:BB:49:C:O2'	36:BB:50:G:H5'	2.17	0.45
36:BB:61:G:C2	36:BB:62:C:C2	3.05	0.45
38:BD:241:PRO:O	38:BD:242:ARG:HB2	2.16	0.45
38:BD:39:LYS:HB2	38:BD:62:TYR:CB	2.47	0.45
39:BE:1:MET:H2	39:BE:84:PHE:HB2	1.82	0.45
40:BF:139:PHE:O	40:BF:142:TRP:HB3	2.17	0.45
40:BF:77:ASP:C	40:BF:79:GLY:N	2.71	0.45
41:BG:133:LEU:C	41:BG:133:LEU:CD1	2.86	0.45
41:BG:77:ILE:C	41:BG:79:ASN:N	2.70	0.45
42:BH:146:ALA:O	42:BH:148:ILE:N	2.50	0.45
42:BH:38:SER:C	42:BH:40:GLU:N	2.71	0.45
42:BH:39:PRO:O	42:BH:42:ARG:HG3	2.17	0.45
42:BH:89:ILE:HG12	42:BH:90:LYS:N	2.32	0.45
43:BI:87:LYS:CE	43:BI:121:LYS:HG2	2.47	0.45
43:BI:97:ILE:CD1	43:BI:116:LEU:HD22	2.46	0.45
44:BN:30:ILE:O	44:BN:30:ILE:HG22	2.16	0.45
44:BN:35:ARG:HD3	44:BN:37:LYS:HD3	1.99	0.45
44:BN:38:HIS:O	44:BN:40:PRO:HD3	2.16	0.45
35:BA:2392:A:H1'	46:BP:60:MET:HE3	1.99	0.45
46:BP:59:LEU:CA	46:BP:61:ARG:CZ	2.90	0.45
48:BR:41:ALA:C	48:BR:43:GLU:N	2.68	0.45
49:BS:25:ARG:HA	49:BS:87:PHE:O	2.17	0.45
50:BT:6:LEU:O	50:BT:6:LEU:HD23	2.17	0.45
51:BU:108:GLU:OE1	51:BU:112:ARG:NH2	2.46	0.45
54:BX:35:THR:HG23	54:BX:36:LYS:N	2.32	0.45
55:BY:31:LEU:CD1	55:BY:34:LYS:N	2.77	0.45
56:BZ:27:VAL:HG13	56:BZ:27:VAL:O	2.16	0.45
1:CA:1230:C:C2'	1:CA:1231:G:H5'	2.46	0.45
1:CA:1246:C:H2'	1:CA:1247:U:H6	1.81	0.45
1:CA:1255:G:H5'	3:CC:26:LYS:NZ	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1437:C:H2'	1:CA:1438:G:O5'	2.17	0.45
1:CA:1446:U:C2	1:CA:1457:G:N1	2.85	0.45
1:CA:1490:C:H2'	1:CA:1491:G:O4'	2.17	0.45
1:CA:299:G:C6	1:CA:300:A:C6	3.05	0.45
1:CA:473:G:H5''	16:CP:81:ARG:NE	2.30	0.45
1:CA:691:G:O2'	1:CA:692:U:H5'	2.17	0.45
1:CA:712:A:O2'	1:CA:713:G:H5'	2.17	0.45
3:CC:27:LYS:NZ	3:CC:27:LYS:O	2.50	0.45
4:CD:154:ASN:HB2	4:CD:159:ARG:HH21	1.82	0.45
5:CE:136:MET:HB3	5:CE:140:ARG:HH21	1.81	0.45
5:CE:76:ILE:HG23	5:CE:77:PRO:CD	2.46	0.45
7:CG:104:LEU:H	7:CG:104:LEU:HD22	1.82	0.45
7:CG:92:SER:CB	7:CG:94:ARG:HE	2.30	0.45
8:CH:31:PHE:HA	8:CH:34:GLU:HG2	1.98	0.45
1:CA:1118:C:OP1	9:CI:104:ARG:HD3	2.16	0.45
10:CJ:48:THR:CB	10:CJ:62:HIS:HB3	2.47	0.45
11:CK:23:ALA:O	11:CK:87:THR:N	2.48	0.45
14:CN:24:CYS:HB3	14:CN:27:CYS:O	2.16	0.45
14:CN:26:ARG:HB2	14:CN:39:LEU:CD2	2.47	0.45
17:CQ:60:ILE:HG12	17:CQ:61:GLU:N	2.31	0.45
20:CT:12:ALA:C	20:CT:14:LYS:H	2.20	0.45
25:CY:161:ILE:O	25:CY:162:GLN:C	2.56	0.45
31:D5:25:LEU:HD12	53:DW:23:LEU:HD11	1.98	0.45
33:D7:39:ARG:HG2	33:D7:39:ARG:HH11	1.82	0.45
35:DA:1018:C:H2'	35:DA:1019:U:H6	1.82	0.45
35:DA:1364:G:H2'	35:DA:1364:G:N3	2.32	0.45
35:DA:1450:G:H2'	35:DA:1450(A):C:H6	1.81	0.45
35:DA:1586:A:N1	35:DA:1587:A:C4	2.85	0.45
35:DA:1668:A:C5	35:DA:1674:G:C5	3.04	0.45
35:DA:729:G:C4	35:DA:1775:U:C2	3.05	0.45
35:DA:1937:A:H2'	35:DA:1938:A:H5'	1.99	0.45
35:DA:2023:G:O2'	35:DA:2024:G:H5'	2.17	0.45
35:DA:2245:U:H5'	35:DA:2246:G:C5'	2.41	0.45
35:DA:2359:C:H2'	35:DA:2360:A:O4'	2.16	0.45
35:DA:2560:C:C2'	35:DA:2561:A:H5'	2.46	0.45
35:DA:2660:A:H2'	35:DA:2661:G:O4'	2.16	0.45
35:DA:2654:A:N1	35:DA:2665:A:H5''	2.32	0.45
35:DA:2713:A:C2'	35:DA:2714:G:H5'	2.47	0.45
35:DA:2828:C:H2'	35:DA:2829:C:H6	1.80	0.45
35:DA:35:G:O2'	35:DA:36:G:H5'	2.17	0.45
35:DA:523:C:C4	35:DA:524:U:C5	3.05	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:764:A:C2	35:DA:781:A:C2	3.05	0.45
36:DB:97:G:C2	36:DB:98:G:C8	3.05	0.45
39:DE:65:GLY:O	39:DE:66:HIS:C	2.55	0.45
40:DF:102:PRO:O	40:DF:103:LYS:C	2.55	0.45
40:DF:107:LYS:O	40:DF:108:LYS:C	2.55	0.45
40:DF:114:VAL:HG12	40:DF:192:LEU:CD1	2.47	0.45
40:DF:132:VAL:CG2	40:DF:133:ASN:H	2.05	0.45
47:DQ:34:LEU:HG	47:DQ:103:MET:HB2	1.99	0.45
35:DA:1652:A:H62	48:DR:11:ASN:ND2	2.14	0.45
49:DS:66:ALA:O	49:DS:67:ARG:HB2	2.17	0.45
50:DT:100:TYR:CD2	50:DT:103:ARG:CZ	2.95	0.45
50:DT:109:GLU:O	50:DT:110:ILE:C	2.54	0.45
50:DT:11:GLU:N	50:DT:11:GLU:OE1	2.47	0.45
50:DT:83:ILE:O	50:DT:84:GLN:C	2.55	0.45
51:DU:43:GLY:HA3	52:DV:76:LYS:HB2	1.99	0.45
53:DW:62:HIS:O	53:DW:64:MET:HG3	2.16	0.45
54:DX:26:TYR:HD1	54:DX:26:TYR:H	1.62	0.45
55:DY:11:ASP:OD1	55:DY:12:THR:N	2.50	0.45
55:DY:22:GLY:O	55:DY:23:ARG:CG	2.59	0.45
56:DZ:156:LYS:O	56:DZ:156:LYS:HG2	2.17	0.45
1:AA:1307:U:C4'	13:AM:109:THR:HG21	2.47	0.44
1:AA:1422:G:H1	1:AA:1479:C:N4	2.14	0.44
1:AA:1521:G:H2'	1:AA:1522:U:C6	2.51	0.44
1:AA:190:U:O2'	1:AA:191:G:H5'	2.16	0.44
1:AA:313:A:H2'	1:AA:314:C:H6	1.76	0.44
1:AA:318:G:H2'	1:AA:319:G:H8	1.81	0.44
1:AA:756:C:H2'	1:AA:757:U:O4'	2.17	0.44
1:AA:790:A:C6	1:AA:791:G:C6	3.04	0.44
1:AA:839:U:O2	1:AA:839:U:O4'	2.33	0.44
2:AB:188:ALA:HB1	2:AB:192:SER:HB2	1.99	0.44
3:AC:134:ILE:HD11	3:AC:153:VAL:CG2	2.46	0.44
4:AD:162:LEU:O	4:AD:165:MET:HB2	2.17	0.44
8:AH:87:SER:OG	8:AH:92:ARG:HA	2.17	0.44
10:AJ:22:LYS:O	10:AJ:24:VAL:N	2.41	0.44
12:AL:76:ASN:HD21	12:AL:108:ALA:HB2	1.82	0.44
13:AM:89:GLY:C	13:AM:90:LEU:O	2.55	0.44
16:AP:48:TRP:CE3	16:AP:49:LEU:N	2.83	0.44
17:AQ:47:PRO:HG2	17:AQ:48:GLU:CD	2.37	0.44
19:AS:40:ILE:HG21	19:AS:62:ILE:HD11	1.99	0.44
20:AT:51:GLU:O	20:AT:52:ALA:C	2.55	0.44
20:AT:63:ILE:O	20:AT:64:ASP:C	2.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AU:21:TYR:HD1	21:AU:21:TYR:N	2.14	0.44
23:AW:44:A:H2'	23:AW:45:A:C8	2.53	0.44
27:B1:56:GLN:HG3	27:B1:57:GLU:N	2.31	0.44
32:B6:26:ASN:O	32:B6:27:LYS:HD3	2.16	0.44
33:B7:28:ARG:O	33:B7:29:LYS:C	2.56	0.44
34:B8:29:LYS:HZ1	34:B8:44:LYS:HB3	1.82	0.44
34:B8:32:LEU:HD23	34:B8:35:GLN:C	2.38	0.44
35:BA:1254:A:H5'	35:BA:1255:U:C5'	2.47	0.44
35:BA:1301:A:O2'	35:BA:1302:A:C2'	2.59	0.44
35:BA:1353:A:C5	35:BA:1378:A:C5	3.05	0.44
35:BA:1388:G:H1	35:BA:1399:C:N4	2.15	0.44
35:BA:1578:U:OP2	35:BA:1578:U:C6	2.69	0.44
35:BA:151:C:N4	35:BA:175:G:H1	2.15	0.44
35:BA:1906:G:C8	35:BA:1929:G:C4	3.05	0.44
35:BA:2020:A:C5	35:BA:2022:U:C5	3.05	0.44
35:BA:2330:G:N2	35:BA:2386:C:C2	2.85	0.44
35:BA:2498:C:C2'	35:BA:2499:C:H5'	2.47	0.44
35:BA:2740:A:C6	35:BA:2764:A:C8	3.05	0.44
35:BA:447:A:C2	35:BA:473:G:C8	3.05	0.44
35:BA:608:A:N3	35:BA:608:A:H2'	2.32	0.44
35:BA:639:U:C4	35:BA:640:C:N4	2.86	0.44
35:BA:785:G:C4	35:BA:786:C:C5	3.05	0.44
35:BA:845:G:C8	35:BA:845:G:OP2	2.69	0.44
36:BB:64:C:O2'	36:BB:65:C:H5'	2.17	0.44
37:BC:65:PRO:HG2	37:BC:189:ILE:CA	2.48	0.44
37:BC:86:ALA:HA	37:BC:89:ALA:HB2	1.98	0.44
40:BF:102:PRO:O	40:BF:104:LYS:N	2.50	0.44
40:BF:148:LEU:CD2	40:BF:191:ARG:HD3	2.47	0.44
40:BF:53:THR:CG2	40:BF:56:GLU:OE1	2.65	0.44
40:BF:5:ALA:C	40:BF:6:VAL:HG22	2.38	0.44
41:BG:111:LEU:HB3	41:BG:117:PHE:CE2	2.51	0.44
41:BG:135:LEU:HG	41:BG:136:ARG:N	2.28	0.44
41:BG:169:ALA:CA	41:BG:173:LEU:HD23	2.47	0.44
41:BG:173:LEU:N	41:BG:173:LEU:CD2	2.72	0.44
42:BH:92:ILE:CG2	42:BH:93:GLY:N	2.73	0.44
44:BN:44:PRO:O	44:BN:46:VAL:N	2.49	0.44
45:BO:113:LYS:HA	45:BO:116:SER:OG	2.16	0.44
47:BQ:50:ALA:O	47:BQ:54:MET:CB	2.65	0.44
48:BR:97:VAL:HG13	48:BR:114:VAL:HG23	1.98	0.44
48:BR:28:LEU:HD12	48:BR:29:LEU:CD1	2.47	0.44
48:BR:53:HIS:CA	48:BR:56:LYS:HB2	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:76:ALA:HB3	50:BT:75:ILE:CG1	2.46	0.44
52:BV:78:LYS:CD	52:BV:79:VAL:N	2.68	0.44
53:BW:43:GLY:O	53:BW:44:ALA:C	2.55	0.44
53:BW:96:ILE:HG13	53:BW:97:LYS:H	1.82	0.44
54:BX:53:LYS:HZ2	54:BX:55:ASN:HD21	1.63	0.44
55:BY:18:GLY:O	55:BY:20:TYR:N	2.51	0.44
55:BY:77:PRO:O	55:BY:78:ALA:CB	2.64	0.44
55:BY:86:ARG:NH2	55:BY:95:LYS:NZ	2.63	0.44
1:CA:1106:G:H5''	3:CC:172:ARG:HG2	1.99	0.44
1:CA:1232:U:H2'	1:CA:1233:G:O4'	2.17	0.44
1:CA:124:G:C6	1:CA:125:U:C4	3.05	0.44
1:CA:402:G:O2'	1:CA:403:C:H5'	2.17	0.44
1:CA:511:C:O2'	1:CA:512:U:H6	2.00	0.44
1:CA:651:C:O2'	1:CA:652:U:H5'	2.17	0.44
4:CD:192:GLU:C	4:CD:194:LEU:N	2.70	0.44
4:CD:29:PRO:C	4:CD:30:LYS:HG2	2.38	0.44
4:CD:65:ARG:NH1	4:CD:72:GLU:CA	2.80	0.44
5:CE:36:ASP:CG	5:CE:37:ARG:N	2.70	0.44
5:CE:60:TYR:HE2	5:CE:64:ARG:NE	2.15	0.44
6:CF:10:LEU:HD13	6:CF:61:LEU:CD1	2.46	0.44
6:CF:69:GLU:CD	6:CF:69:GLU:N	2.70	0.44
7:CG:47:CYS:CA	7:CG:58:PRO:HG3	2.47	0.44
7:CG:64:GLN:O	7:CG:66:VAL:N	2.50	0.44
9:CI:58:ARG:HD3	9:CI:59:PHE:CE1	2.51	0.44
1:CA:1125:U:O4	10:CJ:5:ARG:HD3	2.17	0.44
12:CL:119:LYS:O	12:CL:120:TYR:CB	2.62	0.44
12:CL:38:THR:CG2	12:CL:57:LYS:HB2	2.47	0.44
12:CL:84:LEU:HB3	12:CL:101:VAL:HB	2.00	0.44
12:CL:97:ARG:C	12:CL:98:TYR:CD1	2.90	0.44
18:CR:53:ARG:HH22	18:CR:59:SER:C	2.20	0.44
19:CS:40:ILE:HD13	19:CS:62:ILE:CD1	2.47	0.44
19:CS:6:LYS:HG2	19:CS:7:LYS:NZ	2.32	0.44
27:D1:46:LEU:HD22	27:D1:48:LYS:CE	2.47	0.44
29:D3:16:PRO:HB2	29:D3:18:ASP:OD1	2.16	0.44
35:DA:1264:G:H2'	35:DA:1265:A:C8	2.51	0.44
35:DA:1289:C:O2'	35:DA:1290:C:H5'	2.16	0.44
35:DA:1431:U:O2'	35:DA:1432:C:H5'	2.16	0.44
35:DA:1444:G:C2	35:DA:1548:C:C2	3.05	0.44
35:DA:1831:G:O2'	35:DA:1832:C:H5'	2.16	0.44
35:DA:1923:U:H2'	35:DA:1924:C:H6	1.82	0.44
35:DA:1991:U:H2'	35:DA:1992:G:H5''	1.96	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2000:G:C2	35:DA:2001:A:C8	3.05	0.44
27:D1:41:ARG:NH2	35:DA:205:G:C6	2.85	0.44
35:DA:2295:C:O2'	35:DA:2296:U:H5'	2.17	0.44
35:DA:2313:C:H2'	35:DA:2314:C:C6	2.52	0.44
35:DA:2056:G:C8	35:DA:2577:A:C6	3.05	0.44
35:DA:2636:U:H2'	35:DA:2637:U:H6	1.81	0.44
35:DA:472:A:H2'	35:DA:473:G:H5'	1.99	0.44
35:DA:563:G:C6	35:DA:564:C:N4	2.85	0.44
35:DA:577:G:H2'	35:DA:578:A:H8	1.80	0.44
35:DA:717:G:H2'	35:DA:718:A:O4'	2.16	0.44
35:DA:809:G:H2'	35:DA:810:U:C6	2.52	0.44
35:DA:977:G:C6	35:DA:987:G:C6	3.04	0.44
36:DB:74:U:H3'	36:DB:75:G:H5''	1.94	0.44
37:DC:105:ASP:O	37:DC:106:ALA:HB2	2.16	0.44
37:DC:70:LYS:C	37:DC:72:VAL:H	2.19	0.44
39:DE:14:ILE:O	39:DE:20:ALA:HA	2.18	0.44
39:DE:38:THR:C	39:DE:40:GLU:N	2.69	0.44
40:DF:123:LEU:HD12	40:DF:124:LEU:N	2.31	0.44
40:DF:45:ARG:HG3	40:DF:46:ARG:N	2.32	0.44
40:DF:63:LYS:HZ3	40:DF:67:GLN:HB3	1.80	0.44
44:DN:22:THR:N	44:DN:61:ARG:HB2	2.32	0.44
45:DO:107:ARG:NH1	50:DT:36:GLU:N	2.61	0.44
45:DO:114:ILE:N	45:DO:114:ILE:CD1	2.71	0.44
46:DP:47:ASP:CG	46:DP:49:ARG:HB3	2.37	0.44
35:DA:1242:A:N1	46:DP:8:PRO:CG	2.79	0.44
49:DS:26:LEU:HG	49:DS:39:ILE:HD11	1.99	0.44
51:DU:21:ALA:O	51:DU:22:LYS:C	2.55	0.44
53:DW:8:ARG:HG3	53:DW:8:ARG:NH1	2.31	0.44
56:DZ:145:GLU:HA	56:DZ:145:GLU:OE1	2.17	0.44
1:AA:1419:G:N2	1:AA:1482:G:N9	2.66	0.44
1:AA:1464:G:C2	1:AA:1465:C:C5	3.05	0.44
1:AA:1473:A:O2'	1:AA:1474:G:H5'	2.15	0.44
1:AA:189(K):U:H2'	1:AA:189(L):G:C8	2.52	0.44
1:AA:184:G:O4'	1:AA:224:C:H4'	2.16	0.44
1:AA:402:G:C6	1:AA:403:C:C4	3.04	0.44
1:AA:524:G:H2'	1:AA:525:C:H6	1.79	0.44
1:AA:662:G:O2'	1:AA:836:G:C5'	2.65	0.44
1:AA:824:C:H2'	1:AA:825:G:H8	1.82	0.44
1:AA:76:C:N4	1:AA:93:G:H1	2.14	0.44
3:AC:190:ARG:O	3:AC:191:THR:O	2.35	0.44
3:AC:99:VAL:HG23	3:AC:99:VAL:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:19:LEU:HD23	6:AF:19:LEU:O	2.16	0.44
7:AG:107:ALA:O	7:AG:110:GLN:HB2	2.16	0.44
7:AG:15:ASP:OD2	7:AG:44:TYR:OH	2.35	0.44
7:AG:23:VAL:HG12	7:AG:27:ILE:HD11	1.98	0.44
11:AK:62:GLN:C	11:AK:64:ALA:N	2.69	0.44
12:AL:7:ILE:O	12:AL:8:ASN:C	2.56	0.44
15:AO:56:LEU:HA	15:AO:59:MET:CE	2.47	0.44
17:AQ:11:VAL:HG23	17:AQ:20:THR:HB	1.99	0.44
17:AQ:56:VAL:O	17:AQ:76:LEU:HD12	2.17	0.44
19:AS:10:PHE:CD1	19:AS:10:PHE:N	2.85	0.44
19:AS:78:ARG:HH11	19:AS:78:ARG:HG3	1.80	0.44
23:AW:68:C:O2'	23:AW:69:C:H5'	2.17	0.44
25:AY:170:ALA:O	25:AY:172:ALA:N	2.50	0.44
25:AY:34:ASN:ND2	25:AY:36:ALA:HB3	2.32	0.44
26:B0:51:VAL:CG2	26:B0:81:VAL:HG23	2.48	0.44
27:B1:68:PRO:O	27:B1:69:LYS:C	2.54	0.44
29:B3:4:LEU:C	29:B3:4:LEU:HD23	2.37	0.44
31:B5:42:PRO:C	31:B5:43:HIS:CD2	2.91	0.44
35:BA:814:C:H1'	35:BA:1225:G:N2	2.32	0.44
35:BA:1305:C:C2'	35:BA:1306:C:H5'	2.47	0.44
35:BA:1456:G:C4	35:BA:1457:A:N7	2.85	0.44
35:BA:147:U:H2'	35:BA:148:C:C6	2.52	0.44
35:BA:1525:G:H2'	35:BA:1526:G:C8	2.52	0.44
35:BA:154:G:H1	35:BA:172:C:N4	2.08	0.44
35:BA:120:U:O4	35:BA:177:G:C8	2.70	0.44
35:BA:17:G:H4'	51:BU:25:TRP:HH2	1.77	0.44
35:BA:182:A:N3	35:BA:433:C:O2'	2.46	0.44
35:BA:1945:G:C6	35:BA:1946:U:O4	2.71	0.44
35:BA:1992:G:N1	35:BA:1997:G:N1	2.65	0.44
35:BA:2034:U:C2'	35:BA:2035:G:H5'	2.47	0.44
27:B1:40:ARG:NH1	35:BA:2081:C:O2'	2.50	0.44
35:BA:2245:U:H5'	35:BA:2246:G:C5'	2.41	0.44
35:BA:806:C:O2	35:BA:2444:G:O2'	2.34	0.44
35:BA:253:C:C2'	35:BA:254:G:H5'	2.47	0.44
35:BA:2654:A:N1	35:BA:2665:A:H5''	2.32	0.44
35:BA:2713:A:H3'	35:BA:2714:G:H5'	1.99	0.44
35:BA:2820:A:HO2'	35:BA:2821:A:P	2.40	0.44
35:BA:304:G:C2	35:BA:305:U:C2	3.06	0.44
35:BA:39:C:C4	35:BA:40:C:N4	2.85	0.44
33:B7:16:HIS:ND1	35:BA:465:G:H4'	2.32	0.44
35:BA:523:C:C5	35:BA:524:U:H5	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:543:C:C6	35:BA:547:A:C8	3.04	0.44
35:BA:897:C:O2'	35:BA:898:C:H5'	2.17	0.44
35:BA:93:G:H2'	35:BA:94:C:O4'	2.18	0.44
35:BA:968:G:H2'	35:BA:969:U:C6	2.51	0.44
36:BB:61:G:H2'	36:BB:62:C:H6	1.77	0.44
37:BC:70:LYS:C	37:BC:72:VAL:H	2.20	0.44
38:BD:34:VAL:HG22	38:BD:35:LYS:NZ	2.33	0.44
39:BE:179:GLU:O	39:BE:180:ASN:CB	2.65	0.44
39:BE:78:LEU:C	39:BE:79:ARG:HD2	2.37	0.44
40:BF:39:TRP:CG	40:BF:101:LEU:HB2	2.52	0.44
40:BF:64:ILE:HG12	40:BF:65:TRP:NE1	2.33	0.44
41:BG:170:ARG:NH1	41:BG:170:ARG:HG3	2.33	0.44
42:BH:144:VAL:CG1	42:BH:144:VAL:O	2.65	0.44
42:BH:61:HIS:C	42:BH:63:SER:N	2.71	0.44
42:BH:88:LEU:O	42:BH:89:ILE:CG2	2.65	0.44
44:BN:51:PHE:CZ	44:BN:119:ARG:HD2	2.52	0.44
46:BP:71:VAL:HG22	46:BP:72:PRO:CG	2.47	0.44
47:BQ:125:LEU:HB3	47:BQ:127:ILE:CD1	2.47	0.44
48:BR:17:ARG:NH1	48:BR:17:ARG:HG2	2.28	0.44
49:BS:106:ARG:O	49:BS:106:ARG:HD2	2.17	0.44
49:BS:38:GLN:CG	49:BS:39:ILE:N	2.78	0.44
49:BS:77:ALA:O	49:BS:80:LEU:N	2.45	0.44
50:BT:102:ILE:O	50:BT:103:ARG:C	2.56	0.44
50:BT:75:ILE:CD1	50:BT:75:ILE:N	2.77	0.44
52:BV:75:PHE:HE1	52:BV:89:GLN:HB2	1.83	0.44
53:BW:14:PRO:O	53:BW:16:LYS:N	2.50	0.44
55:BY:28:LYS:CB	55:BY:39:VAL:H	2.30	0.44
1:CA:1065:U:H5'	1:CA:1190:G:N2	2.12	0.44
1:CA:1440:C:N4	1:CA:1441:G:C2	2.85	0.44
1:CA:356:A:H1'	1:CA:368:U:O2'	2.18	0.44
1:CA:66:G:H2'	1:CA:66:G:N3	2.31	0.44
1:CA:756:C:O2'	1:CA:757:U:H5'	2.18	0.44
1:CA:831:U:H2'	1:CA:832:C:C5	2.53	0.44
1:CA:922:G:N3	1:CA:1398:A:C2	2.76	0.44
1:CA:922:G:C6	1:CA:923:A:C6	3.05	0.44
2:CB:180:LEU:C	2:CB:182:ILE:H	2.20	0.44
2:CB:75:LYS:CA	2:CB:75:LYS:HE3	2.48	0.44
3:CC:181:ASN:HD21	3:CC:204:LEU:HB2	1.82	0.44
4:CD:12:CYS:SG	4:CD:18:LYS:HG2	2.57	0.44
5:CE:147:ASP:HA	5:CE:150:ARG:CB	2.43	0.44
11:CK:73:MET:SD	11:CK:103:LEU:CD1	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CK:67:ASP:HA	11:CK:70:LYS:HB3	1.98	0.44
13:CM:89:GLY:O	13:CM:90:LEU:O	2.34	0.44
15:CO:36:ILE:HG22	15:CO:37:ASN:N	2.32	0.44
16:CP:55:ARG:O	16:CP:56:ALA:C	2.56	0.44
18:CR:33:ASP:OD2	18:CR:34:TYR:N	2.49	0.44
19:CS:43:GLU:C	19:CS:45:VAL:N	2.64	0.44
19:CS:33:THR:CG2	19:CS:51:VAL:HG22	2.48	0.44
20:CT:26:ASN:HD22	20:CT:27:LYS:N	2.14	0.44
25:CY:150:SER:N	25:CY:153:GLU:OE1	2.36	0.44
25:CY:15:GLN:N	25:CY:168:PHE:HZ	2.16	0.44
31:D5:15:ARG:HA	31:D5:18:ALA:HB2	1.98	0.44
31:D5:23:HIS:O	31:D5:24:ALA:C	2.53	0.44
32:D6:33:LYS:CA	32:D6:33:LYS:HE2	2.40	0.44
35:DA:1131:G:N2	35:DA:1132:A:C5	2.85	0.44
35:DA:1175:U:O3'	35:DA:1176:G:H3'	2.18	0.44
35:DA:1204:A:N1	35:DA:1241:A:H2	2.15	0.44
35:DA:1478:G:HO2'	35:DA:1558:A:H2	1.66	0.44
35:DA:1526:G:H2'	35:DA:1527:G:O4'	2.18	0.44
35:DA:1464:C:H1'	35:DA:1528(A):A:N3	2.32	0.44
35:DA:1591:G:H2'	35:DA:1592:C:H5'	1.99	0.44
35:DA:1655:A:O2'	39:DE:115:GLY:HA3	2.18	0.44
35:DA:1790:C:H2'	35:DA:1791:A:N7	2.33	0.44
35:DA:2101:G:C6	35:DA:2102:U:C6	3.05	0.44
35:DA:1889:A:N1	35:DA:2234:G:H1'	2.32	0.44
35:DA:2426:A:C3'	35:DA:2427:C:C5'	2.84	0.44
35:DA:2464:C:O2'	35:DA:2465:C:O5'	2.35	0.44
35:DA:2531:A:H2	35:DA:2658:C:O2	2.00	0.44
35:DA:2714:G:C5	35:DA:2715:C:C4	3.05	0.44
35:DA:2775:A:O2'	35:DA:2776:A:C5'	2.64	0.44
35:DA:532:A:H4'	35:DA:533:G:O4'	2.17	0.44
35:DA:641:C:H2'	35:DA:642:G:O4'	2.17	0.44
35:DA:926:A:H5'	35:DA:926:A:C8	2.51	0.44
36:DB:75:G:N7	36:DB:76:G:N7	2.65	0.44
35:DA:1843:C:H1'	38:DD:255:LYS:HZ3	1.82	0.44
38:DD:94:LEU:HD13	38:DD:94:LEU:C	2.37	0.44
40:DF:167:ALA:O	40:DF:168:ARG:C	2.55	0.44
41:DG:131:TYR:O	41:DG:159:VAL:HG21	2.17	0.44
41:DG:135:LEU:HB3	41:DG:155:MET:HE3	1.99	0.44
41:DG:23:PHE:HZ	41:DG:167:GLU:HB3	1.82	0.44
36:DB:57:A:H1'	41:DG:30:GLU:CA	2.47	0.44
35:DA:2313:C:H4'	41:DG:40:ASN:ND2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:17:VAL:HG12	42:DH:19:VAL:HG23	1.99	0.44
45:DO:101:PRO:C	45:DO:102:VAL:HG22	2.38	0.44
45:DO:85:VAL:CG1	45:DO:86:ILE:N	2.80	0.44
46:DP:138:LEU:HD23	46:DP:142:GLY:HA3	1.98	0.44
46:DP:29:LYS:N	46:DP:29:LYS:HD2	2.32	0.44
47:DQ:134:ARG:HH12	56:DZ:119:GLU:CD	2.20	0.44
47:DQ:73:PRO:HG3	47:DQ:93:TYR:CE2	2.53	0.44
49:DS:25:ARG:HA	49:DS:87:PHE:O	2.18	0.44
50:DT:53:ARG:HG2	50:DT:53:ARG:O	2.16	0.44
50:DT:32:TYR:HD2	50:DT:81:PRO:CB	2.30	0.44
50:DT:92:GLY:C	50:DT:94:ALA:H	2.18	0.44
52:DV:73:SER:OG	52:DV:89:GLN:O	2.29	0.44
54:DX:32:PRO:O	54:DX:75:ASP:OD2	2.35	0.44
54:DX:84:ALA:O	54:DX:86:GLY:N	2.50	0.44
56:DZ:165:VAL:HG12	56:DZ:166:SER:N	2.31	0.44
1:AA:1011:G:C2'	1:AA:1012:U:H5'	2.47	0.44
1:AA:101:A:H2'	1:AA:102:G:C8	2.51	0.44
1:AA:1051:C:H2'	1:AA:1052:U:H6	1.83	0.44
1:AA:116:A:H2'	1:AA:117:G:O4'	2.17	0.44
1:AA:1434:A:H2'	1:AA:1435:G:C8	2.52	0.44
1:AA:66:G:N2	1:AA:172:A:C2	2.85	0.44
1:AA:243:A:C2	1:AA:245:C:C2	3.05	0.44
1:AA:429:U:H4'	1:AA:430:A:O5'	2.14	0.44
1:AA:373:A:C8	1:AA:482:A:C8	3.05	0.44
1:AA:794:A:O2'	1:AA:795:C:H5'	2.18	0.44
1:AA:835:U:O2'	1:AA:836:G:H5'	2.18	0.44
1:AA:935:A:H2'	1:AA:936:C:H6	1.83	0.44
2:AB:142:LEU:HD23	2:AB:142:LEU:C	2.37	0.44
3:AC:92:ALA:C	3:AC:94:LEU:N	2.71	0.44
4:AD:114:ARG:O	4:AD:117:ALA:HB3	2.17	0.44
4:AD:192:GLU:O	4:AD:194:LEU:N	2.51	0.44
5:AE:103:GLY:C	5:AE:106:PRO:HD2	2.37	0.44
6:AF:21:LEU:CA	6:AF:24:GLU:HG2	2.46	0.44
6:AF:33:TYR:HD1	6:AF:75:LEU:CG	2.24	0.44
7:AG:11:GLN:HG3	7:AG:12:LEU:N	2.31	0.44
7:AG:148:ASN:C	7:AG:150:ALA:N	2.70	0.44
1:AA:688:G:H5'	11:AK:46:GLY:C	2.38	0.44
12:AL:41:ARG:HG2	12:AL:42:THR:H	1.83	0.44
12:AL:54:LYS:HD2	12:AL:54:LYS:H	1.79	0.44
13:AM:13:LYS:O	13:AM:45:VAL:HG23	2.17	0.44
13:AM:33:ALA:HA	13:AM:59:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AP:73:LEU:CD2	16:AP:73:LEU:N	2.79	0.44
18:AR:29:PHE:HD2	18:AR:29:PHE:H	1.66	0.44
18:AR:53:ARG:HH12	18:AR:59:SER:CA	2.26	0.44
19:AS:33:THR:HG22	19:AS:49:ILE:HG22	2.00	0.44
20:AT:83:ARG:HA	20:AT:86:ARG:HD3	1.98	0.44
35:BA:1042:G:H3'	35:BA:1043:C:O4'	2.18	0.44
35:BA:1174:A:OP1	35:BA:1175:U:H5''	2.17	0.44
35:BA:1180:C:C2'	35:BA:1181:C:H5'	2.46	0.44
35:BA:1231:G:H2'	35:BA:1232:G:H8	1.81	0.44
35:BA:1298:C:H3'	35:BA:1299:G:C8	2.49	0.44
35:BA:1448:G:H5'	35:BA:1449:A:OP1	2.16	0.44
35:BA:1807:G:C2	35:BA:1811:G:O6	2.69	0.44
35:BA:1860:G:H2'	35:BA:1861:G:C8	2.52	0.44
35:BA:2008:C:H2'	35:BA:2009:G:H8	1.83	0.44
35:BA:2075:U:C4	35:BA:2238:G:C6	3.06	0.44
35:BA:234:C:O5'	35:BA:234:C:H6	2.00	0.44
35:BA:2352:A:C2'	35:BA:2353:G:H5'	2.47	0.44
35:BA:2476:A:N3	35:BA:2476:A:H2'	2.31	0.44
35:BA:2617:C:O2'	35:BA:2618:G:H5'	2.17	0.44
35:BA:2712(A):A:OP2	35:BA:2714:G:OP2	2.35	0.44
35:BA:2763:G:C8	35:BA:2763:G:H5'	2.52	0.44
35:BA:2892:A:N6	35:BA:2893:G:N2	2.65	0.44
35:BA:296:C:C2'	35:BA:297:C:H5'	2.48	0.44
35:BA:481:G:O2'	35:BA:482:A:P	2.74	0.44
35:BA:884:C:H4'	35:BA:892:G:C8	2.52	0.44
35:BA:908:C:O2'	35:BA:909:A:H5'	2.17	0.44
35:BA:990:A:O5'	35:BA:991:C:OP2	2.36	0.44
37:BC:75:LEU:HD23	37:BC:75:LEU:C	2.38	0.44
39:BE:87:GLU:O	39:BE:88:GLY:O	2.35	0.44
42:BH:30:LYS:HZ1	42:BH:81:GLU:HA	1.82	0.44
44:BN:96:GLU:HG2	44:BN:97:ARG:N	2.32	0.44
45:BO:43:VAL:HG21	45:BO:52:VAL:CG1	2.47	0.44
45:BO:87:ILE:HD13	45:BO:92:GLU:C	2.38	0.44
50:BT:51:ARG:O	50:BT:61:PHE:HA	2.17	0.44
52:BV:18:LEU:CD2	52:BV:19:LYS:N	2.70	0.44
52:BV:20:LEU:N	52:BV:20:LEU:HD12	2.32	0.44
53:BW:88:ARG:HB2	53:BW:93:ALA:HA	1.99	0.44
55:BY:46:LYS:C	55:BY:47:LYS:HD2	2.38	0.44
56:BZ:11:GLU:HB2	56:BZ:13:GLU:OE2	2.16	0.44
1:CA:101:A:H2'	1:CA:102:G:C8	2.49	0.44
1:CA:1030(B):C:H2'	1:CA:1030(C):G:H5'	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1293:G:O2'	1:CA:1294:G:P	2.76	0.44
1:CA:1300:G:HO2'	1:CA:1301:U:P	2.41	0.44
1:CA:1304:G:N7	1:CA:1305:G:C6	2.85	0.44
1:CA:1365:G:H2'	1:CA:1366:C:O4'	2.17	0.44
1:CA:1504:G:OP1	1:CA:1507:A:H4'	2.17	0.44
1:CA:16:A:C2	1:CA:17:U:C6	3.05	0.44
1:CA:341:C:H2'	1:CA:342:C:H6	1.80	0.44
1:CA:434:U:H2'	1:CA:435:C:C1'	2.47	0.44
1:CA:445:G:N1	1:CA:490:G:C6	2.85	0.44
1:CA:524:G:H2'	1:CA:525:C:H6	1.81	0.44
1:CA:799:G:H2'	1:CA:800:G:H5'	1.99	0.44
1:CA:881:G:P	12:CL:12:ARG:HH22	2.40	0.44
2:CB:115:LEU:CB	2:CB:145:LEU:HD11	2.46	0.44
3:CC:8:ILE:C	3:CC:10:PHE:N	2.70	0.44
3:CC:95:THR:HG22	3:CC:97:LYS:HB2	1.98	0.44
5:CE:110:LEU:HB3	5:CE:115:VAL:HG21	1.99	0.44
5:CE:72:GLN:OE1	5:CE:77:PRO:HA	2.18	0.44
6:CF:37:VAL:HG13	6:CF:65:VAL:HG11	1.99	0.44
7:CG:27:ILE:HD11	7:CG:43:PHE:CE2	2.51	0.44
8:CH:35:ILE:N	8:CH:35:ILE:HD13	2.31	0.44
9:CI:49:PRO:HG3	9:CI:78:LYS:HG2	1.98	0.44
11:CK:64:ALA:O	11:CK:66:LEU:N	2.51	0.44
11:CK:23:ALA:CB	11:CK:91:ARG:HG2	2.47	0.44
13:CM:23:TYR:HB3	13:CM:67:GLU:HA	2.00	0.44
13:CM:73:GLU:HG2	13:CM:77:ASN:HD21	1.83	0.44
14:CN:47:LEU:HB3	14:CN:53:LEU:HG	2.00	0.44
14:CN:47:LEU:O	14:CN:51:GLY:N	2.51	0.44
17:CQ:64:PRO:C	17:CQ:65:ILE:HD12	2.37	0.44
18:CR:26:LEU:HD21	18:CR:42:ARG:NH1	2.32	0.44
19:CS:22:LEU:HD21	19:CS:28:LYS:N	2.33	0.44
20:CT:43:LEU:HA	20:CT:46:GLU:HB3	2.00	0.44
20:CT:26:ASN:CB	20:CT:71:THR:HG23	2.42	0.44
1:CA:261:U:OP2	20:CT:79:ARG:NH2	2.50	0.44
25:CY:160:GLU:C	25:CY:164:ILE:HD13	2.38	0.44
27:D1:14:VAL:O	27:D1:15:ALA:O	2.34	0.44
28:D2:26:ARG:NH2	54:DX:6:ASP:CA	2.75	0.44
35:DA:1051:G:C2	35:DA:1052:C:N4	2.86	0.44
35:DA:121:G:H4'	35:DA:149:A:H5'	1.99	0.44
35:DA:1381:G:O2'	35:DA:1382:G:H5'	2.17	0.44
35:DA:1388:G:O2'	35:DA:1389:G:H5'	2.17	0.44
35:DA:1425:G:H2'	35:DA:1426:G:C8	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:149:A:H2'	35:DA:150:C:H6	1.83	0.44
35:DA:151:C:N4	35:DA:175:G:H1	2.14	0.44
35:DA:1831:G:H2'	35:DA:1832:C:C6	2.52	0.44
35:DA:2027:G:C6	35:DA:2028:U:C4	3.06	0.44
35:DA:2070:G:H2'	35:DA:2071:A:C8	2.52	0.44
35:DA:2088:G:H2'	35:DA:2089:U:O4'	2.17	0.44
35:DA:2388:A:H5'	35:DA:2389:G:OP2	2.18	0.44
35:DA:2498:C:C2'	35:DA:2499:C:H5'	2.46	0.44
35:DA:2727:G:C4	35:DA:2728:U:C5	3.05	0.44
35:DA:2881:C:C4	35:DA:2882:A:N7	2.86	0.44
35:DA:348:G:H2'	35:DA:349:G:H8	1.82	0.44
35:DA:389:G:C8	35:DA:2413:G:H4'	2.53	0.44
35:DA:579:G:C6	35:DA:580:C:N4	2.86	0.44
35:DA:60:G:C5	35:DA:63:U:C4	3.05	0.44
35:DA:824:A:H2'	35:DA:825:C:C6	2.52	0.44
35:DA:926:A:C4	35:DA:927:G:C8	3.06	0.44
35:DA:966:G:H2'	35:DA:967:C:H6	1.83	0.44
35:DA:976:C:H2'	35:DA:977:G:C8	2.53	0.44
38:DD:218:ARG:HG3	38:DD:218:ARG:NH1	2.33	0.44
39:DE:101:ARG:HH11	39:DE:169:ASN:HD22	1.66	0.44
39:DE:4:ILE:HD11	39:DE:29:GLY:H	1.82	0.44
40:DF:160:ASN:HB3	40:DF:163:VAL:CG2	2.46	0.44
40:DF:160:ASN:CB	40:DF:163:VAL:HG23	2.48	0.44
40:DF:178:PRO:HG2	40:DF:179:GLU:HG3	1.98	0.44
40:DF:20:LEU:HD12	40:DF:199:TRP:HH2	1.81	0.44
40:DF:34:TRP:O	40:DF:35:GLU:C	2.55	0.44
42:DH:121:ILE:CG2	42:DH:133:VAL:HG11	2.47	0.44
42:DH:19:VAL:HG22	42:DH:24:VAL:HG13	1.99	0.44
42:DH:54:ARG:HB2	42:DH:61:HIS:HD2	1.83	0.44
43:DI:94:ALA:HB1	43:DI:114:LEU:HD12	1.99	0.44
43:DI:68:LEU:HD23	43:DI:136:VAL:HG11	1.99	0.44
43:DI:93:THR:O	43:DI:96:ASP:OD2	2.35	0.44
46:DP:47:ASP:HB3	46:DP:48:PRO:CA	2.47	0.44
48:DR:24:GLN:CB	48:DR:44:LEU:HD21	2.41	0.44
49:DS:106:ARG:HE	49:DS:106:ARG:HB3	1.51	0.44
49:DS:31:SER:CB	49:DS:34:HIS:O	2.66	0.44
49:DS:38:GLN:HG3	49:DS:49:VAL:O	2.17	0.44
49:DS:53:SER:OG	49:DS:54:LEU:N	2.47	0.44
50:DT:78:LEU:O	50:DT:78:LEU:HD23	2.18	0.44
39:DE:11:MET:N	50:DT:8:LYS:HE3	2.33	0.44
51:DU:111:GLU:C	51:DU:113:ALA:N	2.69	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:118:GLN:HG3	56:DZ:120:ILE:CD1	2.47	0.44
56:DZ:28:MET:CE	56:DZ:37:VAL:HG21	2.48	0.44
1:AA:1030(B):C:H2'	1:AA:1030(C):G:H5'	1.98	0.44
1:AA:1135:U:H2'	1:AA:1137:C:O2	2.18	0.44
1:AA:1348:U:H2'	1:AA:1349:A:H8	1.83	0.44
1:AA:1372:U:OP1	9:AI:71:SER:HB3	2.16	0.44
1:AA:622:A:C8	1:AA:623:C:C6	3.05	0.44
1:AA:658:G:H1'	15:AO:22:THR:CB	2.41	0.44
1:AA:971:G:H5''	1:AA:972:C:H5''	1.99	0.44
3:AC:14:ILE:HG12	3:AC:15:THR:N	2.32	0.44
4:AD:100:ARG:NH2	4:AD:137:SER:HA	2.31	0.44
5:AE:82:VAL:HG11	5:AE:134:ALA:O	2.18	0.44
6:AF:69:GLU:CD	6:AF:69:GLU:N	2.71	0.44
8:AH:35:ILE:HG22	8:AH:111:ILE:HD13	1.99	0.44
8:AH:97:VAL:CG1	8:AH:98:LYS:H	2.18	0.44
1:AA:1118:C:OP1	9:AI:104:ARG:HD3	2.18	0.44
10:AJ:23:ILE:HG23	10:AJ:85:LEU:HD13	1.99	0.44
16:AP:55:ARG:O	16:AP:56:ALA:C	2.55	0.44
25:AY:154:THR:O	25:AY:157:ALA:N	2.51	0.44
25:AY:18:LEU:HD22	25:AY:171:LYS:HB3	2.00	0.44
25:AY:34:ASN:HD21	25:AY:36:ALA:HB3	1.82	0.44
26:B0:84:LEU:HD12	26:B0:84:LEU:C	2.38	0.44
27:B1:43:TYR:HD2	27:B1:45:ASN:HD21	1.63	0.44
32:B6:13:CYS:O	32:B6:21:TYR:HA	2.17	0.44
34:B8:29:LYS:NZ	34:B8:44:LYS:HB2	2.32	0.44
34:B8:59:LYS:HB2	34:B8:59:LYS:NZ	2.27	0.44
35:BA:1278:A:O2'	35:BA:1279:G:H5'	2.17	0.44
35:BA:1396:U:O2	35:BA:1396:U:C2'	2.65	0.44
35:BA:1418:G:H1	35:BA:1579:A:C5'	2.26	0.44
35:BA:1675:C:H2'	35:BA:1676:A:O4'	2.17	0.44
35:BA:1806:C:H42	35:BA:1812:A:N6	2.15	0.44
35:BA:1853:A:N6	35:BA:1889:A:C8	2.86	0.44
31:B5:19:ARG:HG3	35:BA:2046:G:H5'	2.00	0.44
35:BA:2351:G:HO2'	35:BA:2352:A:H8	1.63	0.44
35:BA:2555:U:H2'	35:BA:2556:C:C5'	2.48	0.44
35:BA:2820:A:O2'	35:BA:2821:A:OP1	2.33	0.44
35:BA:2851:A:H2'	35:BA:2852:G:C8	2.52	0.44
35:BA:483:A:H2'	35:BA:484:C:O4'	2.17	0.44
35:BA:644:A:C2'	35:BA:645:C:H5''	2.47	0.44
35:BA:782:A:H8	35:BA:782:A:P	2.40	0.44
35:BA:852:G:C2'	35:BA:853:G:H5'	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:925:C:C2'	35:BA:926:A:C5'	2.84	0.44
35:BA:954:G:N3	35:BA:954:G:H2'	2.33	0.44
36:BB:103:G:O2'	36:BB:104:U:H5'	2.16	0.44
36:BB:115:G:O4'	49:BS:47:THR:HB	2.17	0.44
38:BD:81:ALA:N	38:BD:94:LEU:CD1	2.79	0.44
38:BD:77:ALA:HB1	38:BD:97:TYR:HA	1.99	0.44
39:BE:107:THR:O	39:BE:190:GLY:HA2	2.17	0.44
39:BE:201:THR:HG22	39:BE:203:LYS:HB3	1.99	0.44
40:BF:160:ASN:OD1	40:BF:163:VAL:HG23	2.18	0.44
40:BF:17:ARG:HG3	40:BF:17:ARG:NH1	2.32	0.44
40:BF:32:LEU:HD23	40:BF:33:LEU:N	2.31	0.44
41:BG:138:GLN:NE2	41:BG:152:LEU:HA	2.32	0.44
41:BG:53:LEU:N	41:BG:53:LEU:CD2	2.81	0.44
41:BG:77:ILE:C	41:BG:79:ASN:H	2.21	0.44
43:BI:41:GLU:O	43:BI:44:LEU:HB3	2.17	0.44
47:BQ:134:ARG:CG	47:BQ:135:ASP:N	2.79	0.44
47:BQ:71:ASP:N	47:BQ:94:VAL:O	2.40	0.44
50:BT:119:LYS:O	50:BT:123:GLN:HG2	2.17	0.44
51:BU:58:ARG:HB3	57:BU:201:MG:MG	1.42	0.44
52:BV:43:GLU:HA	52:BV:47:VAL:N	2.33	0.44
52:BV:61:VAL:O	52:BV:62:LEU:HD23	2.18	0.44
52:BV:70:ILE:CG2	52:BV:90:PRO:HB2	2.48	0.44
53:BW:29:LEU:HD21	53:BW:33:ARG:HH21	1.82	0.44
54:BX:31:HIS:O	54:BX:32:PRO:C	2.54	0.44
54:BX:55:ASN:O	54:BX:77:LYS:HB3	2.18	0.44
54:BX:8:ILE:H	54:BX:8:ILE:CD1	2.22	0.44
56:BZ:133:ILE:O	56:BZ:133:ILE:HG22	2.18	0.44
1:CA:63:C:H42	1:CA:104:G:H1	1.64	0.44
1:CA:1057:G:H2'	1:CA:1058:G:C5'	2.47	0.44
1:CA:1150:U:O3'	10:CJ:41:PRO:HA	2.17	0.44
1:CA:1188:A:H2'	1:CA:1189:C:O4'	2.18	0.44
1:CA:1242:C:P	21:CU:10:ARG:HH12	2.40	0.44
1:CA:629:G:H2'	1:CA:630:G:O4'	2.17	0.44
1:CA:64:G:H4'	1:CA:66:G:OP1	2.18	0.44
1:CA:570:G:O6	1:CA:873:A:C2	2.70	0.44
1:CA:935:A:H2'	1:CA:936:C:H6	1.81	0.44
2:CB:12:GLU:O	2:CB:14:GLY:N	2.51	0.44
2:CB:195:ASP:C	2:CB:197:VAL:H	2.19	0.44
3:CC:76:VAL:CG2	3:CC:77:ILE:N	2.81	0.44
4:CD:21:LEU:HD22	4:CD:115:ARG:HG3	1.98	0.44
4:CD:64:LEU:HG	4:CD:65:ARG:N	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:54:THR:OG1	7:CG:56:GLN:HG2	2.18	0.44
8:CH:14:ARG:HB3	8:CH:14:ARG:NH1	2.32	0.44
8:CH:1:MET:HE2	8:CH:2:LEU:H	1.82	0.44
8:CH:45:ILE:HD13	8:CH:62:TYR:O	2.17	0.44
9:CI:11:LYS:C	9:CI:13:ALA:N	2.70	0.44
12:CL:45:PRO:HA	12:CL:93:LEU:CD2	2.47	0.44
13:CM:40:ASN:HA	13:CM:41:PRO:HD3	1.80	0.44
19:CS:41:VAL:CG1	19:CS:44:MET:HB2	2.47	0.44
19:CS:63:THR:HG22	19:CS:66:MET:HE2	1.99	0.44
25:CY:108:GLU:CG	25:CY:111:ARG:NH2	2.80	0.44
25:CY:41:LEU:H	25:CY:41:LEU:CD1	2.23	0.44
27:D1:18:ILE:HG23	27:D1:42:GLN:O	2.16	0.44
29:D3:9:VAL:HG23	29:D3:10:LYS:N	2.33	0.44
33:D7:5:TRP:C	33:D7:6:GLN:NE2	2.70	0.44
35:DA:1239:G:H2'	35:DA:1240:U:O4'	2.17	0.44
35:DA:1284:A:H2'	35:DA:1285:G:O4'	2.17	0.44
35:DA:1309:G:C2'	35:DA:1310:G:H5'	2.48	0.44
35:DA:1358:G:N7	35:DA:1371:G:C6	2.86	0.44
35:DA:142(A):C:C4	35:DA:143:G:N7	2.86	0.44
35:DA:1791:A:OP2	35:DA:1791:A:H8	2.00	0.44
35:DA:2123:G:H21	37:DC:42:GLU:CD	2.20	0.44
35:DA:2262:U:H4'	35:DA:2328:A:H2	1.79	0.44
35:DA:2808:U:H5'	35:DA:2891:G:O6	2.18	0.44
35:DA:306:U:H2'	35:DA:306:U:O2	2.18	0.44
35:DA:60:G:C2	35:DA:74:A:C4	3.05	0.44
35:DA:68:G:N3	35:DA:68:G:H2'	2.32	0.44
35:DA:704:G:N2	35:DA:726:G:C4	2.85	0.44
36:DB:24:G:H4'	36:DB:25:A:N7	2.33	0.44
38:DD:164:GLN:C	38:DD:165:ILE:HD12	2.38	0.44
39:DE:163:GLU:O	39:DE:165:VAL:N	2.51	0.44
39:DE:201:THR:HG22	39:DE:202:LYS:N	2.32	0.44
39:DE:49:LEU:HD23	39:DE:81:ILE:CG1	2.48	0.44
40:DF:148:LEU:HD13	40:DF:154:VAL:HG21	1.99	0.44
40:DF:164:ARG:O	40:DF:166:ALA:N	2.50	0.44
40:DF:123:LEU:HD13	40:DF:192:LEU:HD22	1.99	0.44
40:DF:25:PRO:HG3	40:DF:119:ARG:CA	2.47	0.44
40:DF:5:ALA:C	40:DF:6:VAL:HG22	2.38	0.44
40:DF:70:THR:O	40:DF:72:ARG:N	2.49	0.44
41:DG:102:PHE:CZ	41:DG:141:PHE:HE1	2.35	0.44
42:DH:102:ALA:HB2	42:DH:116:GLU:HA	1.98	0.44
42:DH:38:SER:C	42:DH:40:GLU:N	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:51:PHE:CZ	44:DN:119:ARG:HD2	2.53	0.44
44:DN:23:LEU:O	44:DN:25:ARG:N	2.48	0.44
44:DN:39:ARG:HA	44:DN:40:PRO:HD2	1.86	0.44
44:DN:96:GLU:HG2	44:DN:97:ARG:N	2.32	0.44
46:DP:57:THR:HG22	46:DP:57:THR:O	2.16	0.44
48:DR:19:ALA:O	48:DR:20:LEU:C	2.53	0.44
48:DR:28:LEU:HA	48:DR:34:ILE:CD1	2.47	0.44
48:DR:2:ARG:NE	48:DR:5:LYS:CE	2.81	0.44
50:DT:114:LEU:N	50:DT:114:LEU:HD23	2.31	0.44
51:DU:110:VAL:O	51:DU:113:ALA:HB3	2.17	0.44
52:DV:19:LYS:NZ	52:DV:20:LEU:N	2.60	0.44
35:DA:814:C:OP1	52:DV:84:LYS:HA	2.16	0.44
55:DY:7:VAL:HB	55:DY:8:LYS:NZ	2.32	0.44
56:DZ:150:LEU:HD23	56:DZ:171:ILE:HG13	2.00	0.44
56:DZ:47:VAL:O	56:DZ:48:PHE:C	2.56	0.44
1:AA:1095:U:H5'	1:AA:1109:C:O2	2.17	0.44
1:AA:1129:C:H4'	1:AA:1130:A:H5'	1.98	0.44
1:AA:127:G:OP1	1:AA:635:G:H1'	2.18	0.44
1:AA:1415:G:C2	1:AA:1416:G:C8	3.06	0.44
1:AA:137:C:N4	1:AA:226:G:H1	2.09	0.44
1:AA:59:A:C2	1:AA:331:G:C2	3.06	0.44
1:AA:509:A:C2	1:AA:510:A:C2	3.05	0.44
1:AA:601:C:H42	1:AA:637:G:H1	1.65	0.44
1:AA:611:A:C2'	1:AA:612:C:H5'	2.47	0.44
2:AB:12:GLU:O	2:AB:14:GLY:N	2.51	0.44
5:AE:103:GLY:O	5:AE:106:PRO:CD	2.65	0.44
5:AE:90:VAL:HG21	5:AE:121:LYS:HB3	1.97	0.44
5:AE:12:LEU:HD22	5:AE:12:LEU:C	2.38	0.44
6:AF:83:ASP:C	6:AF:85:VAL:N	2.70	0.44
7:AG:78:ARG:HG3	7:AG:79:ARG:N	2.32	0.44
8:AH:88:LYS:O	8:AH:92:ARG:HD2	2.17	0.44
9:AI:8:GLY:O	9:AI:14:VAL:HG13	2.18	0.44
9:AI:83:ARG:O	9:AI:86:VAL:CG1	2.64	0.44
12:AL:117:ARG:NH2	12:AL:124:LYS:HA	2.32	0.44
12:AL:37:CYS:O	12:AL:79:GLU:O	2.35	0.44
1:AA:1228:C:OP1	13:AM:115:LYS:HG3	2.18	0.44
13:AM:22:ILE:N	13:AM:22:ILE:HD12	2.32	0.44
15:AO:66:LEU:HD12	15:AO:66:LEU:N	2.32	0.44
16:AP:55:ARG:O	16:AP:58:TYR:N	2.50	0.44
6:AF:97:PHE:N	18:AR:30:ASP:OD1	2.49	0.44
19:AS:72:GLY:O	19:AS:74:PHE:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:103:C:OP2	20:AT:14:LYS:HD2	2.17	0.44
21:AU:24:ARG:N	21:AU:24:ARG:HD2	2.32	0.44
25:AY:112:LYS:HG3	25:AY:116:ARG:HD2	1.99	0.44
26:B0:29:GLN:HB2	26:B0:67:VAL:HG23	1.99	0.44
31:B5:16:ARG:CG	31:B5:16:ARG:NH1	2.77	0.44
32:B6:32:ASN:HD22	32:B6:33:LYS:HE3	1.82	0.44
33:B7:31:LEU:HD23	33:B7:42:LEU:HD22	1.98	0.44
34:B8:36:LYS:HE2	34:B8:36:LYS:HB2	1.80	0.44
35:BA:1254:A:H5'	35:BA:1255:U:H5'	1.99	0.44
35:BA:1334:G:O2'	35:BA:1335:U:H5'	2.17	0.44
35:BA:1350:C:C2'	35:BA:1351:C:H5'	2.47	0.44
35:BA:1431:U:C2'	35:BA:1432:C:H5'	2.48	0.44
35:BA:1450:G:H2'	35:BA:1450(A):C:H6	1.83	0.44
35:BA:1467:C:H42	35:BA:1525:G:H1	1.64	0.44
35:BA:1615:C:C5	35:BA:1617:C:C6	3.05	0.44
35:BA:1707:G:C4	35:BA:1708:C:C5	3.05	0.44
35:BA:1989:G:H2'	35:BA:1990:C:H5'	1.99	0.44
35:BA:2073:C:O2'	35:BA:2074:U:H5'	2.17	0.44
35:BA:2463:C:O2'	35:BA:2464:C:H5'	2.17	0.44
35:BA:2752:C:C5	35:BA:2753:A:N7	2.86	0.44
35:BA:2808:U:H5'	35:BA:2891:G:O6	2.17	0.44
35:BA:572:A:H5''	35:BA:573:G:OP2	2.18	0.44
35:BA:727:A:H3'	35:BA:728:G:C8	2.52	0.44
35:BA:968:G:C2	35:BA:969:U:C2	3.05	0.44
35:BA:976:C:H2'	35:BA:977:G:H8	1.83	0.44
38:BD:125:ILE:H	38:BD:125:ILE:HD12	1.77	0.44
39:BE:38:THR:C	39:BE:40:GLU:N	2.70	0.44
39:BE:47:VAL:CG1	39:BE:49:LEU:HD21	2.47	0.44
39:BE:52:LEU:HD23	39:BE:75:VAL:HG22	1.96	0.44
40:BF:102:PRO:O	40:BF:103:LYS:C	2.55	0.44
40:BF:25:PRO:HG3	40:BF:119:ARG:CA	2.48	0.44
40:BF:90:PHE:O	40:BF:91:GLY:C	2.56	0.44
41:BG:173:LEU:O	41:BG:174:GLU:C	2.56	0.44
42:BH:130:ARG:CB	42:BH:130:ARG:HH11	2.30	0.44
44:BN:26:LEU:HG	44:BN:27:ALA:N	2.31	0.44
45:BO:87:ILE:HD13	45:BO:87:ILE:HA	1.87	0.44
45:BO:93:PRO:C	45:BO:95:GLY:N	2.69	0.44
46:BP:48:PRO:HG2	46:BP:49:ARG:H	1.83	0.44
46:BP:48:PRO:O	46:BP:50:ARG:N	2.51	0.44
46:BP:96:THR:HB	46:BP:97:PRO:HD2	2.00	0.44
35:BA:2467:C:H4'	47:BQ:123:HIS:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BR:73:VAL:HG23	48:BR:74:LYS:H	1.82	0.44
53:BW:13:SER:HA	53:BW:14:PRO:HD3	1.79	0.44
56:BZ:128:VAL:HG21	56:BZ:132:ASN:O	2.18	0.44
56:BZ:76:LEU:HD23	56:BZ:76:LEU:N	2.33	0.44
1:CA:1097:C:H2'	1:CA:1098:C:H6	1.81	0.44
1:CA:1168:A:H2'	1:CA:1169:A:H8	1.83	0.44
1:CA:1203:C:H2'	1:CA:1204:A:O4'	2.18	0.44
1:CA:1223:C:OP2	1:CA:1224:G:H8	2.00	0.44
1:CA:1315:U:O2'	1:CA:1316:G:H5'	2.17	0.44
1:CA:1464:G:C2	1:CA:1465:C:C5	3.06	0.44
1:CA:149:A:C2	1:CA:150:C:C4	3.06	0.44
1:CA:1502:A:C2	1:CA:1505:G:N2	2.83	0.44
1:CA:52:G:C6	1:CA:360:A:C2	3.05	0.44
1:CA:50:A:N3	1:CA:52:G:H1'	2.33	0.44
1:CA:573:A:C2	1:CA:574:A:C2	3.05	0.44
1:CA:761:G:H2'	1:CA:762:C:C6	2.53	0.44
1:CA:767:A:H2'	1:CA:768:A:C8	2.51	0.44
1:CA:783:C:C6	1:CA:784:C:H5	2.36	0.44
1:CA:757:U:O2'	1:CA:879:C:H1'	2.17	0.44
2:CB:152:PHE:C	2:CB:152:PHE:CD1	2.91	0.44
2:CB:211:ILE:O	2:CB:212:GLN:C	2.56	0.44
4:CD:98:GLU:C	4:CD:100:ARG:N	2.71	0.44
5:CE:136:MET:O	5:CE:137:GLU:C	2.56	0.44
7:CG:47:CYS:HB3	7:CG:58:PRO:CG	2.48	0.44
8:CH:26:VAL:CG2	8:CH:32:LYS:HZ3	2.27	0.44
9:CI:27:THR:HG23	9:CI:31:GLN:O	2.17	0.44
9:CI:28:VAL:CG1	9:CI:29:ASN:H	2.31	0.44
12:CL:104:VAL:HG12	12:CL:105:TYR:N	2.32	0.44
12:CL:76:ASN:HD21	12:CL:108:ALA:HB2	1.83	0.44
13:CM:16:ASP:OD2	13:CM:16:ASP:N	2.51	0.44
15:CO:60:VAL:O	15:CO:63:ARG:N	2.49	0.44
15:CO:62:GLN:O	15:CO:63:ARG:C	2.55	0.44
16:CP:39:TYR:CD1	16:CP:39:TYR:C	2.90	0.44
17:CQ:9:VAL:O	17:CQ:21:VAL:HG13	2.18	0.44
19:CS:52:TYR:CG	19:CS:53:ASN:N	2.86	0.44
23:CW:52:C:H2'	23:CW:53:G:O4'	2.17	0.44
25:CY:61:PRO:HD2	25:CY:65:THR:O	2.17	0.44
30:D4:13:ARG:O	30:D4:15:ILE:N	2.43	0.44
34:D8:32:LEU:HD11	34:D8:41:ILE:HG22	1.98	0.44
35:DA:1042:G:H3'	35:DA:1043:C:O4'	2.18	0.44
35:DA:1217:C:OP2	51:DU:15:LYS:NZ	2.27	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1267:U:C2'	35:DA:1267:U:O2	2.64	0.44
35:DA:1684:C:O2'	35:DA:1685:C:H5'	2.18	0.44
35:DA:1825:A:H2'	35:DA:1826:G:C8	2.52	0.44
35:DA:18:C:C2	35:DA:19:C:C5	3.06	0.44
35:DA:1955:U:O2'	35:DA:1956:U:H5'	2.17	0.44
35:DA:2034:U:C2'	35:DA:2035:G:H5'	2.47	0.44
35:DA:2282:G:O2'	35:DA:2283:C:OP2	2.24	0.44
35:DA:2475:C:N4	35:DA:2529:G:H22	1.97	0.44
35:DA:1983:C:H4'	35:DA:2606:C:H4'	1.99	0.44
35:DA:2811:G:H5''	35:DA:2811:G:H8	1.82	0.44
35:DA:27:G:C2'	35:DA:28:A:OP2	2.66	0.44
35:DA:601:C:H2'	35:DA:602:G:O4'	2.17	0.44
35:DA:627:A:H8	35:DA:627:A:OP1	2.00	0.44
35:DA:632:A:C2	35:DA:2403:C:H1'	2.52	0.44
35:DA:715:G:H2'	35:DA:716:A:C8	2.53	0.44
35:DA:908:C:O2'	35:DA:909:A:H5'	2.17	0.44
35:DA:942:G:C6	35:DA:943:U:N3	2.85	0.44
35:DA:948:G:H2'	35:DA:949:C:C6	2.53	0.44
36:DB:16:G:H2'	36:DB:17:C:H6	1.83	0.44
36:DB:51:G:H2'	36:DB:52:A:C1'	2.48	0.44
36:DB:78:A:H2'	36:DB:79:C:O4'	2.17	0.44
37:DC:82:LYS:O	37:DC:83:ILE:HD13	2.17	0.44
38:DD:53:PHE:HB3	38:DD:218:ARG:HB2	1.99	0.44
40:DF:103:LYS:HA	40:DF:106:ARG:HE	1.83	0.44
40:DF:192:LEU:C	40:DF:192:LEU:HD23	2.37	0.44
41:DG:56:ALA:O	41:DG:60:LEU:HB2	2.18	0.44
42:DH:92:ILE:O	42:DH:94:TYR:N	2.49	0.44
44:DN:110:GLY:O	44:DN:111:PRO:C	2.55	0.44
44:DN:38:HIS:CG	44:DN:39:ARG:N	2.86	0.44
44:DN:46:VAL:CG2	44:DN:48:MET:HG3	2.48	0.44
46:DP:47:ASP:HB2	46:DP:51:PHE:HD2	1.83	0.44
46:DP:82:GLY:HA2	46:DP:113:LYS:O	2.17	0.44
48:DR:73:VAL:HG23	48:DR:74:LYS:HD3	1.98	0.44
50:DT:78:LEU:HD23	50:DT:79:HIS:ND1	2.33	0.44
51:DU:63:VAL:O	51:DU:64:ARG:C	2.55	0.44
52:DV:72:VAL:O	52:DV:73:SER:HB3	2.17	0.44
52:DV:90:PRO:CD	52:DV:91:TYR:H	2.31	0.44
53:DW:43:GLY:O	53:DW:44:ALA:C	2.56	0.44
56:DZ:53:ILE:CG2	56:DZ:71:VAL:HB	2.47	0.44
1:AA:1111:A:N1	3:AC:177:THR:OG1	2.45	0.44
1:AA:1312:G:N2	1:AA:1326:C:C2	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1441:G:H4'	1:AA:1442:G:C4	2.53	0.44
1:AA:189(B):C:H2'	1:AA:189(C):C:H6	1.82	0.44
1:AA:562:C:O2'	12:AL:15:ARG:HB3	2.18	0.44
1:AA:686:U:H1'	1:AA:687:A:C8	2.52	0.44
1:AA:814:A:C8	1:AA:816:A:C8	3.06	0.44
1:AA:889:A:H5'	1:AA:891:U:O4'	2.17	0.44
2:AB:152:PHE:CD1	2:AB:152:PHE:C	2.90	0.44
2:AB:19:HIS:O	2:AB:20:GLU:O	2.35	0.44
2:AB:24:TRP:CZ3	2:AB:26:PRO:HA	2.53	0.44
4:AD:33:MET:C	4:AD:35:ARG:H	2.20	0.44
5:AE:112:LEU:O	5:AE:114:GLY:N	2.50	0.44
5:AE:139:LEU:H	5:AE:139:LEU:HG	1.48	0.44
5:AE:145:LYS:C	5:AE:148:VAL:HB	2.38	0.44
5:AE:153:LYS:HB3	5:AE:154:GLY:H	1.59	0.44
8:AH:63:LEU:HB3	8:AH:64:LYS:H	1.64	0.44
9:AI:42:ARG:NH2	9:AI:75:ASP:OD1	2.50	0.44
9:AI:18:PHE:HD1	9:AI:62:TYR:CD2	2.35	0.44
11:AK:21:ILE:HG23	11:AK:30:VAL:HG12	2.00	0.44
12:AL:119:LYS:O	12:AL:120:TYR:CB	2.63	0.44
12:AL:45:PRO:HA	12:AL:93:LEU:CD2	2.47	0.44
13:AM:4:ILE:HG22	13:AM:5:ALA:H	1.82	0.44
13:AM:54:VAL:O	13:AM:58:GLU:HG2	2.18	0.44
15:AO:18:PHE:O	15:AO:19:PRO:C	2.55	0.44
15:AO:80:ALA:O	15:AO:81:LEU:C	2.55	0.44
15:AO:87:ILE:O	15:AO:88:ARG:HB2	2.18	0.44
25:AY:144:ALA:O	25:AY:147:LEU:O	2.35	0.44
35:BA:1051:G:C2'	35:BA:1052:C:H5"	2.47	0.44
35:BA:1298:C:C3'	35:BA:1299:G:H8	2.29	0.44
35:BA:1784:A:H4'	35:BA:1785:A:O5'	2.18	0.44
35:BA:1832:C:H2'	35:BA:1833:U:O4'	2.16	0.44
35:BA:1862:G:C2	35:BA:1863:G:C8	3.06	0.44
35:BA:1912:A:H5'	35:BA:1918:A:N1	2.33	0.44
35:BA:1992:G:C2	35:BA:1997:G:C6	3.06	0.44
35:BA:2068:U:N3	35:BA:2430:A:C2	2.80	0.44
35:BA:2569:G:C2	35:BA:2570:G:C8	3.05	0.44
35:BA:2679:A:H2'	35:BA:2680:C:H6	1.82	0.44
35:BA:2831:G:O2'	35:BA:2883:A:H2'	2.18	0.44
35:BA:460:A:C2	35:BA:470:A:C4	3.06	0.44
35:BA:601:C:H2'	35:BA:602:G:O4'	2.18	0.44
35:BA:839:U:H2'	35:BA:840:C:C6	2.53	0.44
35:BA:870:A:N1	35:BA:871:U:C2	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BB:24:G:H4'	36:BB:25:A:H8	1.81	0.44
36:BB:52:A:O2'	36:BB:53:A:C8	2.67	0.44
38:BD:92:ILE:CA	38:BD:107:ALA:HB2	2.48	0.44
35:BA:2579:C:H4'	39:BE:134:ILE:CD1	2.47	0.44
35:BA:323:G:C3'	40:BF:169:ASN:HD21	2.30	0.44
40:BF:117:ARG:NH2	40:BF:187:VAL:HA	2.32	0.44
40:BF:186:ILE:HG23	40:BF:192:LEU:HD12	1.98	0.44
42:BH:160:LYS:O	42:BH:163:TYR:CE1	2.70	0.44
42:BH:35:VAL:O	42:BH:37:VAL:HG23	2.18	0.44
44:BN:41:ASP:O	44:BN:42:TRP:C	2.55	0.44
44:BN:55:VAL:CG1	44:BN:126:PRO:HA	2.48	0.44
44:BN:72:TYR:HB2	44:BN:85:ILE:HB	1.99	0.44
47:BQ:52:VAL:CG1	47:BQ:53:ALA:N	2.67	0.44
49:BS:54:LEU:O	49:BS:56:LEU:N	2.48	0.44
49:BS:72:ALA:C	49:BS:76:LYS:HG2	2.38	0.44
50:BT:30:VAL:CG1	50:BT:44:ASP:HA	2.38	0.44
53:BW:65:LEU:O	53:BW:69:LEU:HG	2.17	0.44
53:BW:80:PRO:O	53:BW:100:THR:HG22	2.17	0.44
54:BX:29:TRP:HE3	54:BX:76:ARG:HB3	1.82	0.44
55:BY:88:LYS:O	55:BY:90:LEU:HD23	2.18	0.44
56:BZ:10:ARG:HG2	56:BZ:11:GLU:N	2.28	0.44
56:BZ:120:ILE:O	56:BZ:121:HIS:HB2	2.18	0.44
56:BZ:24:LEU:HB2	56:BZ:41:LEU:HD23	1.99	0.44
56:BZ:3:TYR:N	56:BZ:57:ILE:HG13	2.33	0.44
56:BZ:69:THR:HG22	56:BZ:90:VAL:HG22	2.00	0.44
1:CA:1307:U:C4'	13:CM:109:THR:HG21	2.48	0.44
1:CA:132:C:N4	1:CA:231:G:N1	2.65	0.44
1:CA:1368:G:H4'	14:CN:61:TRP:HZ2	1.83	0.44
1:CA:1530:G:C4	1:CA:1531:A:C8	3.06	0.44
1:CA:137:C:N4	1:CA:226:G:H1	2.10	0.44
1:CA:346:G:H2'	1:CA:346:G:N3	2.32	0.44
1:CA:39:G:O2'	1:CA:40:C:H5'	2.17	0.44
1:CA:738:C:C2	1:CA:739:C:C5	3.06	0.44
2:CB:236:TYR:HA	2:CB:239:VAL:HG21	1.97	0.44
2:CB:83:MET:HB2	2:CB:84:GLU:H	1.52	0.44
4:CD:64:LEU:O	4:CD:65:ARG:C	2.55	0.44
4:CD:68:TYR:CE2	4:CD:97:LEU:HD22	2.52	0.44
5:CE:20:GLN:O	5:CE:21:ALA:C	2.56	0.44
7:CG:115:ARG:HB2	7:CG:118:VAL:CG2	2.48	0.44
7:CG:145:ALA:C	7:CG:147:ALA:H	2.19	0.44
7:CG:27:ILE:HG23	7:CG:40:ALA:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CK:73:MET:SD	11:CK:103:LEU:HD13	2.57	0.44
16:CP:9:PHE:HE2	16:CP:18:ARG:NE	2.15	0.44
16:CP:29:ASP:N	16:CP:29:ASP:OD2	2.49	0.44
16:CP:67:THR:HG21	16:CP:69:THR:HG23	2.00	0.44
17:CQ:100:LYS:HA	17:CQ:100:LYS:HD3	1.80	0.44
19:CS:12:ASP:HB3	19:CS:15:LEU:CD2	2.45	0.44
19:CS:22:LEU:HD22	19:CS:27:GLU:HB2	2.00	0.44
19:CS:36:ARG:HB3	19:CS:36:ARG:HH11	1.82	0.44
19:CS:78:ARG:CD	19:CS:78:ARG:H	2.31	0.44
21:CU:12:LYS:O	21:CU:22:ARG:NH1	2.51	0.44
25:CY:34:ASN:ND2	25:CY:36:ALA:HB3	2.33	0.44
33:D7:30:VAL:O	33:D7:34:ARG:N	2.42	0.44
34:D8:25:MET:HB2	46:DP:62:LEU:CD1	2.48	0.44
35:DA:1142(A):A:C5	35:DA:1144:G:N7	2.85	0.44
35:DA:1152:C:O2'	35:DA:1153:C:H5'	2.18	0.44
35:DA:812:C:O2	35:DA:1250:G:N1	2.51	0.44
35:DA:139(A):G:N2	54:DX:44:GLU:OE1	2.50	0.44
35:DA:1406:U:H2'	35:DA:1407:C:C6	2.52	0.44
35:DA:1608:A:C5	35:DA:1611:C:C4	3.06	0.44
35:DA:1882:C:H5'	35:DA:1883:G:OP2	2.17	0.44
35:DA:191:A:C2	35:DA:192:C:C2	3.05	0.44
35:DA:2039:C:H2'	35:DA:2040:C:C6	2.52	0.44
35:DA:2194:G:C4	35:DA:2195:C:C5	3.06	0.44
35:DA:2208:A:H1'	35:DA:2219:G:N3	2.32	0.44
35:DA:2420:C:O2'	35:DA:2421:G:H5'	2.18	0.44
35:DA:2778:A:H4'	35:DA:2779:U:OP2	2.17	0.44
35:DA:2793:G:C2	35:DA:2794:C:N3	2.85	0.44
35:DA:2831:G:P	39:DE:58:ARG:NH1	2.90	0.44
35:DA:468:G:C2	35:DA:469:G:H1'	2.53	0.44
35:DA:745:G:P	39:DE:133:LYS:HE3	2.57	0.44
36:DB:74:U:C5	36:DB:75:G:N7	2.85	0.44
37:DC:21:THR:O	37:DC:22:ILE:C	2.55	0.44
38:DD:147:LEU:HB3	38:DD:148:GLU:H	1.59	0.44
38:DD:233:HIS:O	38:DD:234:GLY:C	2.56	0.44
38:DD:247:ALA:HB2	38:DD:253:GLN:HA	2.00	0.44
38:DD:39:LYS:HB2	38:DD:62:TYR:CB	2.48	0.44
39:DE:105:THR:HG22	39:DE:106:GLY:N	2.33	0.44
35:DA:2052:G:P	39:DE:141:ILE:HD11	2.58	0.44
39:DE:16:ARG:O	39:DE:18:ASP:N	2.51	0.44
40:DF:39:TRP:CG	40:DF:101:LEU:HB2	2.50	0.44
40:DF:53:THR:C	40:DF:55:GLY:N	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:117:PHE:CZ	41:DG:179:PRO:HG3	2.53	0.44
35:DA:2312:U:OP1	41:DG:73:ALA:HA	2.18	0.44
43:DI:94:ALA:HA	43:DI:97:ILE:HG13	2.00	0.44
44:DN:56:ASN:HA	44:DN:124:ALA:CA	2.37	0.44
46:DP:85:LEU:HB2	46:DP:120:ALA:HB2	1.99	0.44
46:DP:71:VAL:CB	46:DP:72:PRO:HD3	2.43	0.44
47:DQ:16:ARG:CG	47:DQ:17:LEU:H	2.22	0.44
47:DQ:54:MET:HG3	47:DQ:64:ILE:HD13	1.99	0.44
48:DR:4:LEU:O	48:DR:6:SER:N	2.51	0.44
50:DT:29:ARG:HG3	50:DT:84:GLN:O	2.17	0.44
50:DT:51:ARG:HB3	50:DT:62:THR:OG1	2.18	0.44
50:DT:61:PHE:CE1	50:DT:76:PHE:HD1	2.36	0.44
51:DU:8:VAL:O	51:DU:9:VAL:C	2.56	0.44
52:DV:88:ARG:HG3	52:DV:88:ARG:HH11	1.83	0.44
52:DV:93:GLU:HG2	52:DV:94:LEU:H	1.82	0.44
54:DX:55:ASN:O	54:DX:77:LYS:CB	2.65	0.44
54:DX:60:ARG:CB	54:DX:72:LYS:H	2.31	0.44
54:DX:75:ASP:C	54:DX:76:ARG:HG3	2.38	0.44
55:DY:15:VAL:HG12	55:DY:17:SER:H	1.82	0.44
55:DY:81:LYS:HD3	55:DY:97:ARG:N	2.32	0.44
1:AA:1030(D):A:C2'	1:AA:1031:G:H5'	2.47	0.44
1:AA:1057:G:H2'	1:AA:1058:G:C5'	2.47	0.44
1:AA:1079:G:O3'	5:AE:14:ARG:NH2	2.51	0.44
1:AA:1232:U:H2'	1:AA:1233:G:O4'	2.17	0.44
1:AA:1255:G:H5'	3:AC:26:LYS:NZ	2.32	0.44
1:AA:125:U:H2'	1:AA:126:G:C8	2.53	0.44
1:AA:1300:G:HO2'	1:AA:1301:U:P	2.41	0.44
1:AA:1323:G:H4'	1:AA:1363:C:O2	2.18	0.44
1:AA:132:C:N4	1:AA:231:G:N1	2.65	0.44
1:AA:1346:A:N6	1:AA:1374:A:H3'	2.31	0.44
1:AA:1457:G:C2	1:AA:1458:G:C5	3.05	0.44
1:AA:227:G:C6	1:AA:228:A:C6	3.05	0.44
1:AA:226:G:O2'	1:AA:227:G:H5'	2.18	0.44
1:AA:401:C:O5'	1:AA:401:C:H6	2.01	0.44
1:AA:570:G:C6	1:AA:873:A:N1	2.85	0.44
1:AA:638:G:O2'	1:AA:639:G:H5'	2.18	0.44
1:AA:719:C:O2	18:AR:50:ILE:N	2.36	0.44
1:AA:810:C:H2'	1:AA:811:C:O4'	2.17	0.44
1:AA:858:G:O6	1:AA:869:G:H3'	2.17	0.44
1:AA:961:U:O2'	1:AA:962:C:H5'	2.18	0.44
2:AB:142:LEU:HD23	2:AB:142:LEU:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:213:LEU:HD22	2:AB:214:ILE:HG12	2.00	0.44
5:AE:37:ARG:NH1	5:AE:37:ARG:HG2	2.33	0.44
6:AF:24:GLU:HG3	6:AF:25:ILE:HD13	1.99	0.44
7:AG:104:LEU:O	7:AG:107:ALA:HB3	2.17	0.44
8:AH:125:ARG:HG3	8:AH:125:ARG:HH11	1.83	0.44
9:AI:58:ARG:HD3	9:AI:59:PHE:CE1	2.53	0.44
11:AK:72:ALA:O	11:AK:77:MET:HB2	2.17	0.44
12:AL:27:LEU:O	12:AL:28:LYS:C	2.56	0.44
13:AM:10:PRO:HB2	13:AM:18:ALA:HB1	2.00	0.44
1:AA:973:G:O3'	14:AN:41:ARG:NH2	2.51	0.44
3:AC:37:GLN:NE2	14:AN:52:GLN:OE1	2.50	0.44
18:AR:66:LEU:O	18:AR:69:THR:OG1	2.30	0.44
18:AR:76:LEU:O	18:AR:78:LEU:HG	2.17	0.44
25:AY:66:LEU:HB2	25:AY:101:ILE:HD13	1.99	0.44
25:AY:133:ARG:NH1	25:AY:162:GLN:OE1	2.51	0.44
28:B2:44:LEU:HD12	28:B2:44:LEU:HA	1.76	0.44
29:B3:23:LEU:H	29:B3:23:LEU:HD12	1.83	0.44
34:B8:32:LEU:HD11	34:B8:41:ILE:HG22	1.99	0.44
35:BA:1175:U:C4'	35:BA:1176:G:H3'	2.45	0.44
35:BA:1204:A:N1	35:BA:1241:A:H2	2.16	0.44
35:BA:1298:C:H2'	35:BA:1299:G:H8	1.82	0.44
35:BA:1608:A:H1'	35:BA:1610:A:OP2	2.17	0.44
35:BA:1638:C:OP1	35:BA:2710:C:O2'	2.36	0.44
35:BA:1813:G:N3	38:BD:50:THR:HB	2.33	0.44
35:BA:1910:G:C2	35:BA:1921:G:C4	3.06	0.44
35:BA:2121:G:C6	35:BA:2176:A:N6	2.86	0.44
35:BA:2254:C:C5	35:BA:2255:G:N7	2.86	0.44
35:BA:2299:G:C2	35:BA:2318:G:C8	3.05	0.44
35:BA:2646:C:H2'	35:BA:2647:U:C6	2.52	0.44
35:BA:2727:G:C4	35:BA:2728:U:C5	3.06	0.44
35:BA:2792:G:O6	35:BA:2804:C:N3	2.51	0.44
35:BA:2819:G:H2'	35:BA:2821:A:N7	2.32	0.44
35:BA:49:A:OP1	35:BA:50:U:H3'	2.17	0.44
35:BA:666:G:C5	35:BA:667:U:C5	3.06	0.44
35:BA:737:C:H2'	35:BA:738:G:O5'	2.17	0.44
36:BB:39:A:H2'	36:BB:39:A:N3	2.33	0.44
38:BD:61:LEU:HD12	38:BD:62:TYR:H	1.83	0.44
40:BF:34:TRP:HA	40:BF:37:VAL:CG2	2.48	0.44
41:BG:101:ILE:HG23	41:BG:102:PHE:H	1.83	0.44
41:BG:120:LEU:HD22	41:BG:133:LEU:CD2	2.48	0.44
41:BG:33:ARG:HB2	41:BG:162:THR:OG1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BN:55:VAL:HG11	44:BN:127:ASP:H	1.83	0.44
46:BP:101:VAL:CB	46:BP:107:LYS:HA	2.48	0.44
46:BP:30:THR:O	46:BP:33:ARG:N	2.46	0.44
48:BR:75:LEU:O	48:BR:79:LEU:HB2	2.18	0.44
49:BS:28:VAL:H	49:BS:89:ARG:CB	2.29	0.44
52:BV:36:PRO:HG2	52:BV:60:GLU:CD	2.38	0.44
52:BV:4:ILE:CD1	52:BV:40:LEU:HD11	2.48	0.44
54:BX:35:THR:O	54:BX:36:LYS:C	2.56	0.44
56:BZ:157:LEU:HA	56:BZ:158:PRO:HD2	1.68	0.44
47:BQ:137:TYR:HE2	56:BZ:76:LEU:HD22	1.83	0.44
1:CA:1157:A:H4'	1:CA:1158:C:O5'	2.17	0.44
1:CA:1229:A:H2'	1:CA:1230:C:C6	2.52	0.44
1:CA:126:G:H5'	1:CA:633:G:H22	1.81	0.44
1:CA:1298:C:O4'	1:CA:1299:A:C4	2.71	0.44
1:CA:1457:G:O2'	1:CA:1458:G:H5'	2.17	0.44
1:CA:269:C:H2'	1:CA:270:A:C8	2.52	0.44
1:CA:511:C:C2	1:CA:512:U:C4	3.06	0.44
1:CA:954:G:C5	1:CA:955:U:C4	3.06	0.44
2:CB:164:VAL:O	2:CB:165:VAL:O	2.36	0.44
2:CB:203:GLY:O	2:CB:204:ASN:O	2.36	0.44
2:CB:213:LEU:C	2:CB:213:LEU:CD2	2.86	0.44
2:CB:219:VAL:O	2:CB:222:ILE:HG22	2.17	0.44
3:CC:206:GLU:HB3	3:CC:207:VAL:H	1.58	0.44
3:CC:52:LEU:CD2	3:CC:52:LEU:H	2.21	0.44
4:CD:100:ARG:O	4:CD:101:LEU:C	2.56	0.44
4:CD:33:MET:C	4:CD:35:ARG:H	2.21	0.44
4:CD:56:VAL:C	4:CD:58:LEU:H	2.21	0.44
5:CE:150:ARG:HB2	5:CE:150:ARG:CZ	2.48	0.44
6:CF:100:ASN:CB	18:CR:28:GLU:HA	2.46	0.44
7:CG:102:ARG:HG3	7:CG:106:GLN:NE2	2.33	0.44
7:CG:140:ASP:OD1	7:CG:143:ARG:NH2	2.44	0.44
7:CG:26:PHE:CZ	7:CG:30:ILE:HD11	2.53	0.44
8:CH:86:ILE:HG13	8:CH:133:LEU:HD23	2.00	0.44
10:CJ:22:LYS:O	10:CJ:24:VAL:N	2.42	0.44
11:CK:96:ARG:O	11:CK:99:GLN:HG2	2.17	0.44
15:CO:66:LEU:N	15:CO:66:LEU:HD12	2.32	0.44
19:CS:6:LYS:N	19:CS:6:LYS:CD	2.80	0.44
20:CT:22:ARG:O	20:CT:26:ASN:ND2	2.51	0.44
20:CT:51:GLU:O	20:CT:52:ALA:C	2.56	0.44
23:CW:34:U:O2	23:CW:36:A:C8	2.70	0.44
25:CY:139:LYS:O	25:CY:142:LYS:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:D0:23:VAL:CG1	26:D0:24:LYS:N	2.80	0.44
28:D2:13:ALA:C	28:D2:14:ARG:NE	2.71	0.44
29:D3:23:LEU:O	29:D3:28:LEU:HB2	2.17	0.44
33:D7:34:ARG:C	33:D7:36:GLN:N	2.71	0.44
35:DA:1024:G:P	35:DA:1025:G:H3'	2.58	0.44
35:DA:1353:A:H4'	38:DD:38:LYS:HZ1	1.83	0.44
35:DA:1388:G:H1	35:DA:1399:C:N4	2.14	0.44
35:DA:1778:U:H2'	35:DA:1779:U:H6	1.82	0.44
35:DA:1930:G:H22	35:DA:1968:G:C3'	2.31	0.44
35:DA:1930:G:HO2'	35:DA:1931:U:P	2.40	0.44
35:DA:2121:G:C6	35:DA:2176:A:N6	2.86	0.44
35:DA:2228:G:P	38:DD:263:ARG:HH21	2.40	0.44
35:DA:250:G:H2'	35:DA:251:A:H8	1.83	0.44
35:DA:2512:C:H4'	39:DE:122:PHE:CE2	2.53	0.44
35:DA:1783:A:C2	35:DA:2587:A:C5	3.06	0.44
35:DA:2650:U:H2'	35:DA:2651:C:H6	1.83	0.44
35:DA:2748:A:C2	35:DA:2757:A:C4	3.05	0.44
35:DA:2795:G:O6	35:DA:2801(A):A:C2	2.71	0.44
35:DA:2880:C:H4'	48:DR:90:ARG:NH1	2.33	0.44
31:D5:31:VAL:HG11	35:DA:2886:G:H21	1.83	0.44
35:DA:374:A:C2'	35:DA:375:C:H5'	2.47	0.44
35:DA:483:A:H2'	35:DA:484:C:O4'	2.17	0.44
35:DA:542:C:H2'	35:DA:542:C:O2	2.18	0.44
35:DA:559:G:H2'	35:DA:560:C:H6	1.82	0.44
35:DA:644:A:C2'	35:DA:645:C:H5''	2.48	0.44
35:DA:729:G:N3	35:DA:729:G:H3'	2.32	0.44
38:DD:43:ARG:HB2	38:DD:54:ARG:HB2	1.98	0.44
38:DD:92:ILE:C	38:DD:92:ILE:HD12	2.37	0.44
39:DE:137:HIS:CB	39:DE:138:PRO:HD2	2.36	0.44
39:DE:4:ILE:CD1	39:DE:28:ALA:HB1	2.48	0.44
40:DF:43:LYS:HG3	40:DF:44:ARG:N	2.33	0.44
41:DG:25:TYR:OH	41:DG:168:GLU:HG3	2.18	0.44
35:DA:1035:U:H5''	42:DH:59:ARG:NH1	2.33	0.44
43:DI:42:SER:C	43:DI:44:LEU:H	2.21	0.44
44:DN:17:ASP:O	44:DN:19:GLU:HG3	2.18	0.44
44:DN:44:PRO:O	44:DN:46:VAL:N	2.51	0.44
47:DQ:52:VAL:O	47:DQ:53:ALA:C	2.55	0.44
51:DU:57:PHE:O	51:DU:58:ARG:C	2.55	0.44
52:DV:1:MET:CE	52:DV:46:VAL:HG23	2.48	0.44
54:DX:33:LYS:O	54:DX:34:ALA:C	2.54	0.44
35:DA:1341:U:H5'	54:DX:57:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:DX:8:ILE:H	54:DX:8:ILE:CD1	2.25	0.44
55:DY:28:LYS:NZ	55:DY:37:VAL:HG12	2.28	0.44
1:AA:1009:G:O2'	1:AA:1010:G:H5'	2.18	0.44
1:AA:100:C:H2'	1:AA:101:A:C8	2.53	0.44
1:AA:1049:U:OP1	14:AN:3:ARG:NH1	2.50	0.44
1:AA:949:A:C2	1:AA:1233:G:N3	2.86	0.44
1:AA:1243:C:OP2	21:AU:10:ARG:NH1	2.50	0.44
1:AA:1410:G:N2	1:AA:1491:G:H1'	2.33	0.44
1:AA:583:A:H2'	1:AA:584:G:C8	2.53	0.44
1:AA:708:C:H2'	1:AA:709:G:H8	1.83	0.44
2:AB:77:ALA:HB1	2:AB:211:ILE:HD13	1.99	0.44
3:AC:109:PRO:HA	3:AC:115:LEU:HD12	1.99	0.44
3:AC:11:ARG:O	3:AC:14:ILE:O	2.35	0.44
5:AE:105:VAL:HB	5:AE:106:PRO:HD3	2.00	0.44
6:AF:9:VAL:HG12	6:AF:10:LEU:N	2.33	0.44
7:AG:27:ILE:HD11	7:AG:43:PHE:CD2	2.53	0.44
8:AH:122:ARG:CA	8:AH:125:ARG:HB3	2.43	0.44
8:AH:53:VAL:O	8:AH:54:ASP:CB	2.66	0.44
10:AJ:98:ILE:HG22	10:AJ:99:LYS:N	2.33	0.44
13:AM:4:ILE:HG22	13:AM:5:ALA:N	2.33	0.44
16:AP:23:ASP:OD1	16:AP:24:ALA:N	2.50	0.44
17:AQ:99:SER:O	17:AQ:100:LYS:HE2	2.18	0.44
17:AQ:14:LYS:HB2	17:AQ:14:LYS:HZ3	1.82	0.44
18:AR:37:VAL:O	18:AR:40:LEU:N	2.50	0.44
1:AA:664:G:P	18:AR:64:ARG:HH21	2.41	0.44
19:AS:63:THR:HG22	19:AS:66:MET:CE	2.48	0.44
25:AY:58:VAL:HG22	25:AY:68:VAL:HG13	1.99	0.44
27:B1:52:ARG:NH1	35:BA:2199:A:H5'	2.33	0.44
28:B2:46:GLN:HE21	28:B2:47:ASN:N	2.14	0.44
32:B6:51:GLU:O	32:B6:52:VAL:HB	2.18	0.44
34:B8:2:PRO:O	34:B8:3:LYS:HB2	2.17	0.44
35:BA:1507:A:C2	35:BA:1508:A:H1'	2.53	0.44
35:BA:1517:G:C2'	35:BA:1518:U:H5'	2.48	0.44
35:BA:1720:U:H2'	35:BA:1721:G:C5'	2.48	0.44
35:BA:688:U:H5'	35:BA:1780:A:C2	2.53	0.44
35:BA:2029:G:H2'	35:BA:2031:A:OP2	2.18	0.44
35:BA:2163:C:O2'	35:BA:2164:C:P	2.75	0.44
35:BA:2352:A:H2'	35:BA:2353:G:H5'	2.00	0.44
35:BA:249:C:H5'	35:BA:2394:C:O2'	2.18	0.44
35:BA:2701:C:C3'	35:BA:2702:U:H5''	2.31	0.44
35:BA:422:A:C6	35:BA:423:A:C6	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:426:C:H2'	35:BA:427:U:C5'	2.48	0.44
35:BA:431:U:H6	35:BA:431:U:O5'	2.00	0.44
35:BA:530:G:C5	35:BA:2022:U:H5''	2.52	0.44
35:BA:585:G:H2'	35:BA:586:A:N7	2.32	0.44
35:BA:614:U:C4'	35:BA:614(C):A:H62	2.31	0.44
35:BA:641:C:C2'	35:BA:642:G:H5'	2.48	0.44
35:BA:664:C:H4'	35:BA:941:A:OP1	2.18	0.44
28:B2:52:ASP:OD2	35:BA:72:U:H1'	2.18	0.44
35:BA:761:A:H8	35:BA:761:A:O5'	2.01	0.44
36:BB:103:G:H1'	56:BZ:73:GLN:HE22	1.83	0.44
38:BD:211:ARG:C	38:BD:213:ARG:N	2.69	0.44
40:BF:160:ASN:HB3	40:BF:163:VAL:CG2	2.47	0.44
40:BF:167:ALA:O	40:BF:168:ARG:C	2.56	0.44
40:BF:53:THR:HG23	40:BF:56:GLU:H	1.83	0.44
30:B4:6:HIS:N	41:BG:67:LYS:HE3	2.32	0.44
42:BH:121:ILE:HG22	42:BH:133:VAL:HG11	1.99	0.44
42:BH:13:LYS:CA	42:BH:13:LYS:HE2	2.35	0.44
42:BH:41:MET:HE1	42:BH:55:PRO:HD2	1.99	0.44
42:BH:87:LEU:CD1	42:BH:148:ILE:HG21	2.48	0.44
35:BA:7:G:H4'	44:BN:13:TRP:CZ2	2.52	0.44
44:BN:19:GLU:C	44:BN:21:LYS:H	2.21	0.44
45:BO:47:ILE:HG23	45:BO:48:PRO:N	2.33	0.44
46:BP:113:LYS:HG2	46:BP:115:LEU:CD2	2.48	0.44
51:BU:7:GLY:C	51:BU:8:VAL:CG2	2.86	0.44
52:BV:36:PRO:CG	52:BV:60:GLU:OE1	2.66	0.44
52:BV:46:VAL:HG12	52:BV:47:VAL:N	2.33	0.44
53:BW:75:TYR:HE1	53:BW:104:THR:HB	1.74	0.44
53:BW:19:LEU:O	53:BW:20:VAL:C	2.56	0.44
53:BW:24:ILE:O	53:BW:26:GLY:N	2.50	0.44
47:BQ:138:ASP:OD1	56:BZ:99:TYR:HE1	2.00	0.44
1:CA:949:A:C2	1:CA:1233:G:N3	2.86	0.44
1:CA:1500:A:OP1	1:CA:1505:G:OP1	2.35	0.44
1:CA:241:C:C1'	1:CA:286:G:N2	2.81	0.44
1:CA:385:C:H2'	1:CA:386:C:C6	2.51	0.44
1:CA:433:C:H6	1:CA:433:C:O5'	2.01	0.44
1:CA:452:A:H4'	16:CP:72:ARG:CZ	2.47	0.44
1:CA:677:U:H2'	1:CA:678:U:C6	2.53	0.44
1:CA:725:G:O2'	1:CA:726:C:H5'	2.17	0.44
1:CA:756:C:H2'	1:CA:757:U:O4'	2.18	0.44
3:CC:61:ALA:O	3:CC:62:ASP:HB2	2.18	0.44
4:CD:14:ARG:C	4:CD:16:GLY:N	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CF:37:VAL:HG13	6:CF:65:VAL:HG12	1.99	0.44
6:CF:39:LYS:CG	6:CF:40:VAL:H	2.28	0.44
7:CG:32:ARG:O	7:CG:33:ASP:HB2	2.18	0.44
7:CG:57:GLU:O	7:CG:59:LEU:N	2.49	0.44
8:CH:63:LEU:HB3	8:CH:64:LYS:H	1.64	0.44
11:CK:30:VAL:HG21	11:CK:68:ALA:HB2	2.00	0.44
18:CR:62:GLU:O	18:CR:63:GLN:C	2.56	0.44
20:CT:27:LYS:C	20:CT:27:LYS:HD3	2.38	0.44
23:CW:17:C:H5''	23:CW:18:U:C6	2.53	0.44
25:CY:10:THR:C	25:CY:14:MET:HG3	2.33	0.44
25:CY:25:LEU:C	25:CY:27:GLY:N	2.71	0.44
26:D0:16:SER:CB	35:DA:2261:C:H3'	2.48	0.44
34:D8:56:GLU:HA	34:D8:59:LYS:CE	2.47	0.44
35:DA:1175:U:H4'	35:DA:1176:G:H5'	1.98	0.44
35:DA:1669:A:H5''	35:DA:2550:G:OP1	2.18	0.44
35:DA:1863:G:H1	35:DA:1879:C:H42	1.66	0.44
35:DA:2029:G:C4	35:DA:2031:A:OP2	2.70	0.44
35:DA:1853:A:N1	35:DA:2087:G:H1'	2.33	0.44
35:DA:2128:C:H5'	35:DA:2173:A:H2	1.83	0.44
35:DA:2262:U:H2'	35:DA:2263:C:C5'	2.41	0.44
26:D0:39:ARG:NH2	35:DA:2354:G:N2	2.65	0.44
35:DA:2517:C:C4	35:DA:2542:A:C6	3.06	0.44
35:DA:2660:A:H2'	35:DA:2661:G:C8	2.53	0.44
35:DA:40:C:H2'	35:DA:41:C:C6	2.50	0.44
35:DA:42:G:H2'	35:DA:43:A:H8	1.78	0.44
35:DA:559:G:N2	51:DU:49:HIS:HD2	2.14	0.44
35:DA:559:G:O2'	35:DA:560:C:H5'	2.17	0.44
35:DA:567:A:N1	35:DA:568:U:O2	2.51	0.44
35:DA:695:G:C6	35:DA:696:G:N7	2.86	0.44
35:DA:707:G:H3'	35:DA:708:C:C6	2.53	0.44
35:DA:751:A:H62	35:DA:789:A:N6	2.15	0.44
35:DA:792:G:N3	35:DA:2072:G:H1'	2.33	0.44
36:DB:100:A:C4	36:DB:101:G:C8	3.05	0.44
36:DB:115:G:H2'	36:DB:116:G:C8	2.52	0.44
38:DD:206:LEU:HD23	38:DD:211:ARG:NH1	2.32	0.44
38:DD:268:ARG:HB2	38:DD:268:ARG:CZ	2.47	0.44
38:DD:72:LYS:HB2	38:DD:97:TYR:HE2	1.82	0.44
38:DD:91:ARG:O	38:DD:107:ALA:HB3	2.18	0.44
39:DE:103:ASP:OD1	39:DE:201:THR:HA	2.18	0.44
40:DF:89:VAL:HG12	40:DF:90:PHE:CD1	2.53	0.44
41:DG:109:VAL:O	41:DG:110:ALA:C	2.56	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:131:TYR:CE1	41:DG:133:LEU:HG	2.53	0.44
42:DH:59:ARG:HG2	42:DH:59:ARG:O	2.18	0.44
45:DO:13:ASN:HD22	45:DO:97:ARG:HG2	1.82	0.44
46:DP:48:PRO:CG	46:DP:49:ARG:N	2.81	0.44
46:DP:57:THR:C	46:DP:59:LEU:N	2.70	0.44
47:DQ:70:PRO:HA	47:DQ:94:VAL:C	2.38	0.44
49:DS:56:LEU:O	49:DS:57:LYS:HB3	2.18	0.44
50:DT:32:TYR:HB2	50:DT:33:LYS:H	1.45	0.44
50:DT:65:LYS:HG3	50:DT:66:VAL:N	2.33	0.44
50:DT:79:HIS:O	50:DT:80:SER:HB2	2.18	0.44
50:DT:88:ILE:HG22	50:DT:89:VAL:HG23	2.00	0.44
51:DU:8:VAL:O	51:DU:10:ARG:N	2.51	0.44
52:DV:61:VAL:HG21	52:DV:100:ARG:H	1.82	0.44
1:AA:1438:G:H2'	1:AA:1439:C:H6	1.82	0.44
1:AA:44:G:N2	1:AA:45:U:H1'	2.33	0.44
1:AA:511:C:O2'	1:AA:512:U:H6	2.01	0.44
1:AA:565:U:H2'	1:AA:566:G:H8	1.83	0.44
3:AC:88:ARG:HG2	3:AC:101:LEU:CB	2.47	0.44
4:AD:126:ILE:CG2	4:AD:127:THR:N	2.79	0.44
4:AD:145:GLU:C	4:AD:146:ILE:HD13	2.38	0.44
4:AD:173:TRP:C	4:AD:186:LEU:HB2	2.39	0.44
4:AD:62:GLN:HB3	4:AD:66:ARG:HH22	1.81	0.44
7:AG:115:ARG:HB2	7:AG:118:VAL:CG2	2.48	0.44
9:AI:17:VAL:HG22	9:AI:63:ILE:CG2	2.48	0.44
9:AI:114:TYR:HE1	10:AJ:60:ARG:H	1.58	0.44
11:AK:109:VAL:HG22	18:AR:85:LEU:O	2.18	0.44
12:AL:38:THR:CG2	12:AL:57:LYS:HB2	2.47	0.44
14:AN:26:ARG:HB2	14:AN:39:LEU:CD2	2.47	0.44
14:AN:41:ARG:HG3	14:AN:42:ILE:N	2.31	0.44
15:AO:85:LEU:HB2	15:AO:87:ILE:CD1	2.48	0.44
16:AP:12:LYS:O	16:AP:13:HIS:HB2	2.17	0.44
16:AP:74:LEU:O	16:AP:77:ALA:HB3	2.18	0.44
25:AY:184:LEU:O	25:AY:185:GLY:OXT	2.35	0.44
25:AY:18:LEU:O	25:AY:19:GLU:C	2.56	0.44
27:B1:10:LYS:HA	27:B1:13:ILE:CG2	2.47	0.44
28:B2:14:ARG:NE	28:B2:14:ARG:H	2.15	0.44
29:B3:58:VAL:HG12	29:B3:59:VAL:N	2.33	0.44
34:B8:17:THR:HG22	35:BA:650:C:O2'	2.17	0.44
34:B8:39:LYS:HZ2	34:B8:40:GLU:HA	1.82	0.44
35:BA:1175:U:H4'	35:BA:1176:G:H5'	2.00	0.44
35:BA:1262:A:P	53:BW:99:ARG:HH12	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1484:G:H3'	35:BA:1485:G:C5'	2.46	0.44
35:BA:1615:C:H5	35:BA:1617:C:C5	2.35	0.44
35:BA:2195:C:C2'	35:BA:2196:C:H5'	2.48	0.44
35:BA:2285:C:H6	35:BA:2285:C:H3'	1.83	0.44
35:BA:2795:G:O6	35:BA:2801(A):A:C2	2.71	0.44
35:BA:290:G:O2'	35:BA:291:C:H5'	2.18	0.44
35:BA:718:A:H2'	35:BA:719:C:O4'	2.18	0.44
36:BB:16:G:H2'	36:BB:17:C:C6	2.52	0.44
36:BB:20:C:H2'	36:BB:21:G:C5'	2.38	0.44
40:BF:41:LEU:HA	40:BF:44:ARG:CG	2.48	0.44
41:BG:114:ILE:HD12	41:BG:117:PHE:CD2	2.53	0.44
41:BG:73:ALA:HB2	41:BG:87:PRO:CG	2.44	0.44
42:BH:66:GLY:CA	42:BH:69:ARG:HB2	2.36	0.44
43:BI:112:LYS:C	43:BI:114:LEU:H	2.22	0.44
43:BI:42:SER:C	43:BI:44:LEU:H	2.20	0.44
44:BN:17:ASP:O	44:BN:19:GLU:HG3	2.18	0.44
46:BP:95:VAL:HG23	46:BP:125:VAL:CB	2.46	0.44
47:BQ:70:PRO:HA	47:BQ:94:VAL:C	2.38	0.44
48:BR:72:ASP:HB3	48:BR:75:LEU:HB2	2.00	0.44
51:BU:54:LYS:O	51:BU:58:ARG:HG3	2.18	0.44
53:BW:50:VAL:HG13	53:BW:105:VAL:HG21	1.99	0.44
53:BW:9:TYR:HD2	53:BW:9:TYR:H	1.66	0.44
54:BX:49:VAL:HG13	54:BX:85:PRO:HB3	1.98	0.44
35:BA:1601:G:OP2	54:BX:58:HIS:HD2	2.01	0.44
55:BY:11:ASP:OD1	55:BY:12:THR:N	2.51	0.44
55:BY:15:VAL:HG12	55:BY:16:ALA:N	2.25	0.44
56:BZ:120:ILE:HB	56:BZ:171:ILE:O	2.17	0.44
1:CA:1161:C:H2'	1:CA:1162:C:C6	2.53	0.44
1:CA:1242:C:O5'	21:CU:10:ARG:NH1	2.51	0.44
1:CA:1254:C:H2'	1:CA:1255:G:H8	1.82	0.44
1:CA:125:U:H2'	1:CA:126:G:C8	2.53	0.44
1:CA:1312:G:N2	1:CA:1326:C:C2	2.86	0.44
1:CA:152:A:N6	1:CA:170:U:C2	2.86	0.44
1:CA:339:C:O2'	1:CA:340:U:H5'	2.17	0.44
1:CA:766:A:C5	1:CA:814:A:C2	3.06	0.44
1:CA:828:A:H2'	1:CA:829:G:O4'	2.17	0.44
2:CB:51:LEU:O	2:CB:55:PHE:HD2	2.01	0.44
2:CB:80:ILE:CD1	2:CB:215:LEU:HD12	2.47	0.44
3:CC:35:GLU:OE2	3:CC:95:THR:HG23	2.18	0.44
4:CD:101:LEU:CD2	4:CD:121:VAL:HG13	2.46	0.44
5:CE:41:VAL:CG1	5:CE:113:ALA:HA	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:125:ARG:HG3	8:CH:125:ARG:HH11	1.83	0.44
8:CH:121:ASP:O	8:CH:125:ARG:NE	2.51	0.44
9:CI:3:GLN:HA	9:CI:19:LEU:O	2.17	0.44
9:CI:28:VAL:O	9:CI:30:GLY:N	2.51	0.44
9:CI:42:ARG:NH2	9:CI:75:ASP:OD1	2.51	0.44
9:CI:5:TYR:HE1	9:CI:7:THR:OG1	1.99	0.44
11:CK:127:LYS:O	11:CK:127:LYS:HE2	2.18	0.44
12:CL:21:LYS:CD	12:CL:21:LYS:H	2.09	0.44
13:CM:8:GLU:C	13:CM:9:ILE:HD12	2.38	0.44
14:CN:22:THR:O	14:CN:23:ARG:HB2	2.17	0.44
15:CO:11:VAL:HG13	15:CO:15:PHE:CE1	2.53	0.44
15:CO:6:GLU:N	15:CO:6:GLU:OE1	2.33	0.44
15:CO:83:GLU:O	15:CO:83:GLU:HG2	2.18	0.44
18:CR:22:VAL:HG13	18:CR:25:THR:HB	2.00	0.44
1:CA:1340:A:O2'	22:CV:31:U:H5'	2.18	0.44
27:D1:62:VAL:CG2	27:D1:67:ILE:HA	2.28	0.44
28:D2:32:LEU:CG	28:D2:33:MET:N	2.81	0.44
31:D5:15:ARG:HG3	31:D5:15:ARG:HH11	1.83	0.44
31:D5:40:LYS:NZ	31:D5:46:CYS:H	2.16	0.44
35:DA:1227:G:OP1	51:DU:13:LYS:HE2	2.18	0.44
35:DA:1305:C:C2'	35:DA:1306:C:H5'	2.48	0.44
35:DA:1654:A:O2'	35:DA:1655:A:H5'	2.17	0.44
35:DA:1788:C:C2	35:DA:1789:A:C8	3.06	0.44
35:DA:1802:A:H2'	35:DA:1803:A:C8	2.53	0.44
35:DA:1807:G:C2	35:DA:1811:G:O6	2.71	0.44
35:DA:2004:G:H2'	35:DA:2005:A:C8	2.53	0.44
35:DA:1130:U:O2	35:DA:2025:C:H5''	2.18	0.44
35:DA:2073:C:H2'	35:DA:2074:U:H6	1.83	0.44
35:DA:2584:U:O2	35:DA:2584:U:O4'	2.33	0.44
35:DA:2711:A:C8	35:DA:2714:G:O4'	2.71	0.44
35:DA:2892:A:N6	35:DA:2893:G:N2	2.66	0.44
35:DA:373:U:H1'	35:DA:423:A:N3	2.32	0.44
35:DA:38:A:H2'	35:DA:39:C:H6	1.83	0.44
35:DA:826:U:H2'	35:DA:828:U:O4'	2.18	0.44
35:DA:975(A):G:N3	35:DA:1156:A:H2	2.15	0.44
38:DD:109:ASP:HB2	38:DD:197:GLY:HA2	2.00	0.44
39:DE:140:SER:OG	39:DE:141:ILE:N	2.51	0.44
39:DE:167:VAL:HG22	39:DE:170:LEU:HD11	2.00	0.44
39:DE:33:VAL:HG11	39:DE:89:ASP:N	2.11	0.44
39:DE:51:PHE:CD1	39:DE:51:PHE:C	2.91	0.44
40:DF:192:LEU:HD21	40:DF:194:MET:CE	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DG:16:ARG:HG3	41:DG:16:ARG:NH1	2.32	0.44
41:DG:88:ILE:HG22	41:DG:89:GLY:N	2.33	0.44
41:DG:93:THR:C	41:DG:94:LEU:HD23	2.39	0.44
43:DI:25:TYR:HE2	43:DI:29:TYR:CD2	2.36	0.44
44:DN:104:LYS:C	44:DN:106:MET:H	2.22	0.44
45:DO:17:ARG:O	45:DO:18:LYS:CG	2.66	0.44
45:DO:43:VAL:HG11	45:DO:46:ALA:HB2	1.99	0.44
46:DP:147:LEU:HB2	46:DP:148:LEU:H	1.57	0.44
49:DS:58:LEU:CD2	49:DS:65:VAL:HG13	2.48	0.44
49:DS:72:ALA:C	49:DS:76:LYS:HG2	2.37	0.44
50:DT:121:ILE:CG2	50:DT:122:ASP:N	2.80	0.44
50:DT:29:ARG:HG2	50:DT:86:ILE:N	2.33	0.44
50:DT:57:PHE:CG	50:DT:58:ASN:N	2.85	0.44
50:DT:62:THR:HG22	50:DT:75:ILE:CA	2.23	0.44
51:DU:39:LEU:O	51:DU:40:PHE:C	2.55	0.44
51:DU:74:LEU:HD12	51:DU:74:LEU:O	2.17	0.44
53:DW:65:LEU:O	53:DW:69:LEU:HG	2.18	0.44
53:DW:6:ILE:O	53:DW:6:ILE:HG22	2.17	0.44
55:DY:90:LEU:HD23	55:DY:90:LEU:N	2.33	0.44
56:DZ:33:LEU:CD2	56:DZ:35:ARG:HB2	2.42	0.44
56:DZ:48:PHE:CE2	56:DZ:52:SER:HA	2.53	0.44
1:AA:1086:U:O2'	1:AA:1087:G:H5'	2.18	0.43
1:AA:1299:A:C5	1:AA:1301:U:N3	2.86	0.43
1:AA:1357:A:N7	1:AA:1358:U:C4	2.86	0.43
1:AA:1410:G:C2	1:AA:1491:G:N3	2.86	0.43
1:AA:1442:G:C8	1:AA:1442(B):A:C2	3.06	0.43
1:AA:227:G:C2	1:AA:228:A:C4	3.05	0.43
1:AA:433:C:H6	1:AA:433:C:O5'	2.00	0.43
1:AA:434:U:H2'	1:AA:435:C:C1'	2.48	0.43
1:AA:445:G:N1	1:AA:490:G:C6	2.86	0.43
1:AA:561:U:O2'	1:AA:562:C:P	2.76	0.43
1:AA:735:C:HO2'	1:AA:736:C:H5'	1.80	0.43
1:AA:831:U:H2'	1:AA:832:C:C5	2.53	0.43
2:AB:100:GLY:O	2:AB:104:ASN:C	2.56	0.43
2:AB:212:GLN:HE22	2:AB:216:SER:HB2	1.76	0.43
3:AC:121:ALA:HB1	3:AC:188:LEU:O	2.18	0.43
1:AA:407:G:O2'	4:AD:116:GLN:HG3	2.18	0.43
4:AD:160:GLN:O	4:AD:163:GLU:HB3	2.18	0.43
4:AD:167:GLY:O	4:AD:168:ARG:C	2.56	0.43
4:AD:18:LYS:HE3	4:AD:31:CYS:CB	2.48	0.43
4:AD:74:GLN:C	4:AD:76:ARG:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AE:87:SER:OG	5:AE:125:SER:CB	2.66	0.43
5:AE:147:ASP:O	5:AE:151:LEU:HG	2.18	0.43
1:AA:15:G:H1'	5:AE:19:MET:HG2	1.99	0.43
5:AE:62:ALA:O	5:AE:65:ASN:N	2.47	0.43
5:AE:99:GLY:O	5:AE:117:ASP:HA	2.18	0.43
7:AG:66:VAL:HB	7:AG:67:GLU:OE2	2.18	0.43
9:AI:9:ARG:HA	9:AI:14:VAL:HA	1.99	0.43
12:AL:119:LYS:HD3	12:AL:120:TYR:CE1	2.52	0.43
12:AL:6:THR:H	12:AL:9:GLN:NE2	2.15	0.43
13:AM:91:ARG:CB	13:AM:96:LEU:O	2.63	0.43
13:AM:9:ILE:HD13	41:BG:146:TYR:OH	2.17	0.43
15:AO:33:THR:HG23	15:AO:63:ARG:HH11	1.81	0.43
17:AQ:27:PHE:O	17:AQ:36:ILE:HG13	2.18	0.43
19:AS:22:LEU:HD22	19:AS:27:GLU:HB2	1.99	0.43
20:AT:24:LEU:O	20:AT:27:LYS:CB	2.66	0.43
31:B5:31:VAL:O	31:B5:32:PRO:O	2.36	0.43
35:BA:1042:G:N3	35:BA:1114:G:N2	2.66	0.43
35:BA:1131:G:C2	35:BA:1132:A:N7	2.85	0.43
35:BA:579:G:C2	35:BA:1262:A:C4	3.07	0.43
35:BA:1326:U:C2'	35:BA:1327:C:H5'	2.48	0.43
35:BA:1346:G:H1	35:BA:1600:C:H42	1.65	0.43
35:BA:1767:C:O2	35:BA:1985:G:N2	2.43	0.43
35:BA:1882:C:O2	35:BA:1882:C:H2'	2.18	0.43
35:BA:1889:A:H1'	35:BA:2087:G:O4'	2.18	0.43
35:BA:2052:G:H2'	35:BA:2053:G:C8	2.41	0.43
35:BA:2374:C:O2'	35:BA:2375:G:H5'	2.18	0.43
35:BA:2636:U:H2'	35:BA:2637:U:H6	1.83	0.43
35:BA:2810:A:C2'	39:BE:61:ARG:NH2	2.81	0.43
35:BA:2864:G:H5'	35:BA:2864:G:C8	2.49	0.43
35:BA:2882:A:OP1	48:BR:96:ARG:HD3	2.17	0.43
35:BA:2883:A:C5'	35:BA:2884:U:H5'	2.48	0.43
35:BA:393:C:H2'	35:BA:394:A:C8	2.50	0.43
35:BA:40:C:H2'	35:BA:41:C:C6	2.51	0.43
35:BA:444:C:O2'	35:BA:445:C:H5'	2.18	0.43
35:BA:724:U:H2'	35:BA:725:G:H5'	1.98	0.43
35:BA:855:G:C6	35:BA:856:C:C4	3.06	0.43
38:BD:267:SER:HA	38:BD:270:ILE:CG1	2.48	0.43
38:BD:31:LYS:O	38:BD:32:SER:O	2.35	0.43
38:BD:43:ARG:HB2	38:BD:54:ARG:HB2	2.00	0.43
40:BF:157:VAL:HB	40:BF:194:MET:CB	2.48	0.43
41:BG:108:ASN:C	41:BG:112:PRO:HG2	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:133:LEU:HD12	41:BG:133:LEU:C	2.38	0.43
42:BH:124:GLU:HB2	42:BH:132:ARG:O	2.18	0.43
43:BI:79:ILE:C	43:BI:81:VAL:H	2.21	0.43
43:BI:91:SER:H	43:BI:121:LYS:HE2	1.82	0.43
46:BP:48:PRO:CG	46:BP:49:ARG:H	2.31	0.43
47:BQ:16:ARG:HG2	47:BQ:17:LEU:N	2.25	0.43
48:BR:19:ALA:O	48:BR:20:LEU:C	2.55	0.43
48:BR:2:ARG:HE	48:BR:5:LYS:HZ2	1.66	0.43
35:BA:996:A:C1'	51:BU:92:ARG:HH21	2.31	0.43
56:BZ:19:ARG:HB3	56:BZ:19:ARG:HH11	1.83	0.43
56:BZ:27:VAL:O	56:BZ:88:PHE:N	2.41	0.43
1:CA:1057:G:C2'	1:CA:1058:G:H5'	2.48	0.43
1:CA:1169:A:C2	1:CA:1170:A:C4	3.06	0.43
1:CA:1203:C:OP1	14:CN:3:ARG:CD	2.63	0.43
1:CA:1268:A:H4'	21:CU:20:LYS:H	1.82	0.43
1:CA:1271:G:H5'	1:CA:1314:C:C5'	2.42	0.43
1:CA:1239:A:N6	1:CA:1299:A:H62	2.08	0.43
1:CA:1323:G:H4'	1:CA:1363:C:O2	2.18	0.43
1:CA:1434:A:H2'	1:CA:1435:G:O4'	2.18	0.43
1:CA:1497:G:C6	1:CA:1498:U:O4	2.71	0.43
1:CA:44:G:C2	1:CA:399:G:C2	3.05	0.43
1:CA:561:U:O2'	1:CA:562:C:P	2.76	0.43
1:CA:702:A:C3'	1:CA:703:G:H5'	2.48	0.43
1:CA:783:C:H2'	1:CA:784:C:H6	1.83	0.43
1:CA:824:C:H4'	8:CH:1:MET:H1	1.83	0.43
2:CB:236:TYR:C	2:CB:238:LEU:H	2.20	0.43
7:CG:133:GLY:HA2	7:CG:136:LYS:CG	2.47	0.43
1:CA:933:G:OP2	7:CG:3:ARG:HB3	2.17	0.43
10:CJ:48:THR:HG1	10:CJ:62:HIS:CG	2.36	0.43
1:CA:692:U:H5	11:CK:26:ASN:CG	2.21	0.43
12:CL:47:LYS:HG2	12:CL:48:PRO:HD3	2.00	0.43
14:CN:22:THR:HB	14:CN:33:VAL:HG21	1.99	0.43
17:CQ:29:HIS:ND1	17:CQ:31:LEU:N	2.52	0.43
18:CR:40:LEU:O	18:CR:43:PHE:HD1	2.02	0.43
18:CR:76:LEU:O	18:CR:78:LEU:HG	2.17	0.43
25:CY:10:THR:O	25:CY:12:SER:N	2.51	0.43
25:CY:41:LEU:HD22	25:CY:83:ILE:CD1	2.48	0.43
27:D1:87:PRO:O	27:D1:91:LYS:N	2.40	0.43
28:D2:13:ALA:O	28:D2:14:ARG:O	2.35	0.43
35:DA:118:A:H1'	35:DA:178:G:O4'	2.17	0.43
35:DA:552:G:H1'	35:DA:1220:A:C2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1259:G:H2'	35:DA:1260:G:C8	2.52	0.43
35:DA:1615:C:H5	35:DA:1617:C:C2	2.37	0.43
35:DA:1853:A:H2'	35:DA:1854:A:O4'	2.17	0.43
35:DA:17:G:H2'	35:DA:18:C:C6	2.52	0.43
35:DA:1912:A:H5'	35:DA:1918:A:N1	2.33	0.43
35:DA:1982:C:C2	35:DA:1983:C:C5	3.05	0.43
35:DA:961:C:C5	35:DA:2031:A:C2	3.06	0.43
35:DA:2400:G:H2'	35:DA:2401:U:H6	1.83	0.43
35:DA:2460:U:C2'	35:DA:2461:C:H5'	2.48	0.43
35:DA:261:G:HO2'	35:DA:609:A:H2	1.61	0.43
35:DA:2696:U:H2'	35:DA:2697:G:C8	2.52	0.43
35:DA:2704:C:C2	35:DA:2705:A:C8	3.05	0.43
35:DA:2726:U:HO2'	35:DA:2727:G:C5'	2.31	0.43
35:DA:2831:G:O2'	35:DA:2883:A:H2'	2.17	0.43
35:DA:2832:U:O4	35:DA:2883:A:H5''	2.18	0.43
35:DA:604:G:O6	35:DA:625:G:C6	2.70	0.43
35:DA:604:G:H2'	35:DA:605:C:C6	2.53	0.43
35:DA:685:A:C2	35:DA:787:U:H1'	2.53	0.43
35:DA:724:U:H2'	35:DA:725:G:H5'	2.00	0.43
35:DA:813:U:H2'	35:DA:814:C:H6	1.77	0.43
35:DA:826:U:C2	35:DA:828:U:H1'	2.53	0.43
35:DA:838:C:O2'	35:DA:839:U:H5'	2.17	0.43
35:DA:974:G:C4	35:DA:989:G:C2	3.05	0.43
36:DB:28:C:H42	36:DB:56:G:H1	1.66	0.43
36:DB:64:C:O2'	36:DB:65:C:H5'	2.18	0.43
37:DC:59:ARG:HB2	37:DC:62:VAL:CG2	2.41	0.43
38:DD:139:GLY:H	38:DD:165:ILE:HB	1.82	0.43
38:DD:241:PRO:O	38:DD:242:ARG:HB2	2.18	0.43
38:DD:82:ILE:HA	38:DD:92:ILE:O	2.18	0.43
40:DF:160:ASN:OD1	40:DF:163:VAL:HG23	2.18	0.43
40:DF:182:ASN:O	40:DF:182:ASN:OD1	2.36	0.43
40:DF:21:ALA:C	40:DF:23:ASP:N	2.71	0.43
41:DG:11:TYR:O	41:DG:12:TYR:C	2.56	0.43
41:DG:92:VAL:CG2	41:DG:93:THR:H	2.27	0.43
44:DN:39:ARG:HH11	44:DN:39:ARG:HG3	1.83	0.43
46:DP:131:SER:O	46:DP:135:LEU:N	2.51	0.43
46:DP:16:ARG:O	46:DP:18:ARG:N	2.51	0.43
47:DQ:68:ILE:CD1	47:DQ:68:ILE:N	2.80	0.43
48:DR:81:ASP:O	48:DR:82:GLU:HB2	2.17	0.43
49:DS:26:LEU:CD2	49:DS:28:VAL:HG22	2.47	0.43
51:DU:106:PHE:O	51:DU:107:ALA:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DV:58:VAL:HG12	52:DV:101:GLY:C	2.38	0.43
53:DW:80:PRO:O	53:DW:100:THR:HG22	2.17	0.43
53:DW:67:ASP:N	53:DW:69:LEU:HD11	2.33	0.43
56:DZ:40:ASP:OD1	56:DZ:43:GLU:N	2.34	0.43
1:AA:1054:C:O2'	1:AA:1055:A:H5''	2.18	0.43
1:AA:1342:C:O2'	1:AA:1343:G:H5'	2.17	0.43
1:AA:1502:A:H2	1:AA:1505:G:N1	2.12	0.43
1:AA:192:U:C2	1:AA:193:C:C5	3.06	0.43
1:AA:260:G:H2'	1:AA:261:U:C6	2.53	0.43
1:AA:318:G:C2	1:AA:319:G:C5	3.06	0.43
1:AA:322:C:O2'	1:AA:323:U:H5'	2.18	0.43
1:AA:35:G:C4	1:AA:550:G:N2	2.86	0.43
1:AA:584:G:O2'	1:AA:585:G:H5'	2.18	0.43
1:AA:829:G:H2'	1:AA:830:G:C8	2.53	0.43
1:AA:859:A:O2'	1:AA:860:A:H5'	2.18	0.43
3:AC:27:LYS:O	3:AC:27:LYS:NZ	2.50	0.43
3:AC:61:ALA:O	3:AC:62:ASP:HB2	2.19	0.43
4:AD:138:TYR:C	4:AD:138:TYR:CD2	2.92	0.43
5:AE:102:ALA:CB	5:AE:106:PRO:HG2	2.47	0.43
6:AF:12:PRO:HG2	6:AF:13:ASN:H	1.82	0.43
6:AF:33:TYR:CE1	6:AF:75:LEU:HA	2.53	0.43
6:AF:88:VAL:CG1	6:AF:89:MET:N	2.81	0.43
7:AG:54:THR:OG1	7:AG:56:GLN:HG2	2.18	0.43
8:AH:48:TYR:O	8:AH:49:GLU:HB3	2.19	0.43
1:AA:972:C:C2'	10:AJ:55:LYS:HD3	2.47	0.43
10:AJ:57:LYS:HD2	10:AJ:60:ARG:HH21	1.83	0.43
1:AA:1228:C:O3'	13:AM:116:THR:HA	2.18	0.43
13:AM:19:LEU:O	13:AM:22:ILE:HB	2.18	0.43
13:AM:23:TYR:HB3	13:AM:67:GLU:HA	1.99	0.43
13:AM:73:GLU:HG2	13:AM:77:ASN:HD21	1.82	0.43
15:AO:85:LEU:CD1	15:AO:87:ILE:HD11	2.48	0.43
19:AS:11:VAL:HG22	19:AS:16:LEU:HD11	1.98	0.43
19:AS:6:LYS:HG2	19:AS:7:LYS:NZ	2.32	0.43
20:AT:36:LEU:O	20:AT:37:SER:C	2.56	0.43
25:AY:122:ALA:O	25:AY:126:ARG:NH1	2.51	0.43
25:AY:158:GLU:CD	25:AY:158:GLU:C	2.76	0.43
25:AY:28:LEU:O	25:AY:30:THR:N	2.49	0.43
29:B3:11:SER:OG	29:B3:13:ILE:HG12	2.18	0.43
31:B5:31:VAL:CB	31:B5:32:PRO:HD2	2.34	0.43
33:B7:19:ARG:HH11	33:B7:19:ARG:HG2	1.83	0.43
34:B8:56:GLU:CA	34:B8:59:LYS:NZ	2.80	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:585:G:C5	35:BA:1251:C:C4	3.05	0.43
35:BA:1300:U:H1'	35:BA:1626:G:N2	2.33	0.43
35:BA:1453:U:H4'	35:BA:1455:G:OP1	2.19	0.43
35:BA:1605:C:H2'	35:BA:1606:G:O4'	2.18	0.43
35:BA:2123:G:H21	37:BC:42:GLU:CD	2.21	0.43
35:BA:1889:A:N1	35:BA:2234:G:H1'	2.33	0.43
35:BA:2300:G:H1	35:BA:2316:C:N4	2.10	0.43
35:BA:242:G:N3	35:BA:254:G:C6	2.87	0.43
35:BA:2566:A:N6	45:BO:28:SER:HB2	2.33	0.43
35:BA:2633:G:H5'	35:BA:2811:G:O2'	2.17	0.43
35:BA:2832:U:C5	35:BA:2884:U:H5''	2.54	0.43
35:BA:366:C:C4	35:BA:404:C:C5	3.06	0.43
35:BA:36:G:O2'	35:BA:37:C:H5'	2.18	0.43
35:BA:389:G:H22	46:BP:71:VAL:CG1	2.10	0.43
35:BA:572:A:H2'	35:BA:573:G:O4'	2.18	0.43
35:BA:695:G:C6	35:BA:696:G:N7	2.87	0.43
35:BA:902:C:H2'	35:BA:903:C:H6	1.81	0.43
35:BA:924:C:O2'	35:BA:925:C:H5'	2.18	0.43
36:BB:51:G:H2'	36:BB:52:A:C1'	2.48	0.43
38:BD:69:ARG:C	38:BD:71:ASP:H	2.21	0.43
38:BD:75:ILE:HG21	38:BD:99:ASP:HB2	2.00	0.43
39:BE:169:ASN:OD1	39:BE:201:THR:CG2	2.66	0.43
39:BE:35:GLN:HE22	39:BE:37:ARG:HH21	1.65	0.43
40:BF:114:VAL:HG11	40:BF:202:PHE:CE2	2.47	0.43
40:BF:164:ARG:O	40:BF:165:ARG:C	2.56	0.43
40:BF:3:GLU:HB3	40:BF:20:LEU:O	2.18	0.43
40:BF:22:ALA:HB1	40:BF:26:ALA:CB	2.48	0.43
44:BN:31:ALA:O	44:BN:32:THR:C	2.56	0.43
45:BO:43:VAL:O	45:BO:43:VAL:HG12	2.18	0.43
46:BP:114:ILE:HD11	46:BP:130:PHE:CZ	2.53	0.43
46:BP:47:ASP:OD1	46:BP:50:ARG:HG3	2.19	0.43
46:BP:85:LEU:HB2	46:BP:120:ALA:HB2	2.00	0.43
47:BQ:12:GLN:HE21	47:BQ:72:LYS:HA	1.83	0.43
49:BS:13:ARG:O	49:BS:15:ARG:HD3	2.18	0.43
50:BT:28:VAL:O	50:BT:29:ARG:CD	2.66	0.43
50:BT:48:ILE:CD1	50:BT:48:ILE:N	2.81	0.43
52:BV:29:PRO:C	52:BV:31:ALA:H	2.21	0.43
54:BX:60:ARG:CG	54:BX:74:PRO:HD2	2.33	0.43
1:CA:1050:G:H2'	1:CA:1051:C:H6	1.82	0.43
1:CA:1049:U:O2'	1:CA:1050:G:OP2	2.29	0.43
1:CA:1115:C:H2'	1:CA:1116:C:C6	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1227:A:N7	1:CA:1228:C:C2	2.86	0.43
1:CA:951:G:C6	1:CA:1231:G:C6	3.06	0.43
1:CA:1403:C:C1'	1:CA:1500:A:N1	2.71	0.43
1:CA:515:G:N2	1:CA:537:G:C4	2.87	0.43
2:CB:105:PHE:HA	2:CB:108:ILE:CG2	2.48	0.43
2:CB:114:ARG:HD2	2:CB:118:LEU:HG	1.99	0.43
2:CB:169:LYS:O	2:CB:172:ILE:CD1	2.66	0.43
4:CD:109:GLY:O	4:CD:161:ASN:HB3	2.18	0.43
4:CD:100:ARG:NH2	4:CD:137:SER:HA	2.34	0.43
4:CD:64:LEU:CD2	4:CD:203:VAL:HG21	2.48	0.43
5:CE:147:ASP:CA	5:CE:150:ARG:HH11	2.18	0.43
1:CA:1080:A:C5'	5:CE:16:THR:HG21	2.48	0.43
5:CE:68:GLU:OE2	5:CE:70:PRO:HD3	2.18	0.43
7:CG:30:ILE:HD13	7:CG:105:VAL:HG13	1.99	0.43
7:CG:111:ARG:HH11	7:CG:111:ARG:HB3	1.83	0.43
8:CH:91:ARG:CG	8:CH:91:ARG:NH1	2.79	0.43
10:CJ:32:ALA:N	10:CJ:78:ASN:CG	2.71	0.43
11:CK:29:ILE:CG2	11:CK:44:SER:HB3	2.35	0.43
12:CL:28:LYS:O	12:CL:30:ALA:N	2.51	0.43
19:CS:72:GLY:O	19:CS:74:PHE:N	2.51	0.43
20:CT:71:THR:HB	20:CT:72:LEU:H	1.49	0.43
22:CV:29:G:H2'	22:CV:30:A:H8	1.84	0.43
28:D2:41:ILE:CD1	28:D2:41:ILE:N	2.75	0.43
32:D6:44:ARG:HG2	32:D6:44:ARG:HH11	1.83	0.43
33:D7:16:HIS:CE1	35:DA:465:G:C4'	3.01	0.43
35:DA:99:U:OP1	35:DA:102:G:OP1	2.36	0.43
35:DA:1334:G:O2'	35:DA:1335:U:H5'	2.18	0.43
35:DA:1528(A):A:N7	35:DA:1529:G:C8	2.86	0.43
35:DA:1683:C:H2'	35:DA:1684:C:H6	1.83	0.43
35:DA:1882:C:H2'	35:DA:1882:C:O2	2.17	0.43
35:DA:1935:G:H3'	35:DA:1962:C:N4	2.29	0.43
35:DA:1958:C:O2'	35:DA:1959:G:H5'	2.18	0.43
26:D0:16:SER:HB3	35:DA:2261:C:H3'	2.00	0.43
32:D6:27:LYS:HG3	35:DA:2286:A:OP2	2.16	0.43
35:DA:2287:A:C2	35:DA:2289:G:C8	3.05	0.43
35:DA:2464:C:N4	35:DA:2487:G:N1	2.66	0.43
35:DA:2702:U:OP1	35:DA:2702:U:O4'	2.36	0.43
35:DA:2736:G:H2'	35:DA:2737:G:H8	1.82	0.43
35:DA:462:C:O2'	35:DA:463:G:H5'	2.18	0.43
35:DA:49:A:H4'	35:DA:50:U:H5''	1.99	0.43
35:DA:543:C:C6	35:DA:547:A:C8	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:608:A:H2'	35:DA:608:A:N3	2.33	0.43
35:DA:675:A:O2'	35:DA:676:A:H5'	2.18	0.43
35:DA:78:A:N1	35:DA:109:G:C6	2.87	0.43
35:DA:823:G:N1	35:DA:835:A:C4	2.86	0.43
35:DA:856:C:H4'	35:DA:857:C:OP1	2.18	0.43
35:DA:942:G:H2'	35:DA:943:U:O4'	2.18	0.43
35:DA:947:G:H2'	35:DA:948:G:H8	1.83	0.43
35:DA:986:C:C2'	35:DA:987:G:H5'	2.47	0.43
37:DC:49:ILE:HB	37:DC:50:ASP:H	1.50	0.43
35:DA:1813:G:N3	38:DD:50:THR:HB	2.33	0.43
39:DE:50:GLY:HA3	39:DE:74:PRO:HG3	2.00	0.43
40:DF:114:VAL:CG1	40:DF:202:PHE:HE2	2.31	0.43
41:DG:38:VAL:O	41:DG:38:VAL:HG12	2.17	0.43
45:DO:118:ALA:HA	45:DO:119:PRO:HD2	1.87	0.43
45:DO:67:LYS:O	45:DO:68:GLU:C	2.55	0.43
45:DO:76:ALA:CB	50:DT:75:ILE:HD13	2.47	0.43
45:DO:65:THR:HA	45:DO:82:ASN:HA	2.00	0.43
46:DP:75:ILE:O	46:DP:76:LYS:C	2.56	0.43
47:DQ:57:HIS:ND1	47:DQ:58:PHE:N	2.66	0.43
47:DQ:7:MET:O	47:DQ:8:LYS:HB3	2.18	0.43
48:DR:10:LEU:HB3	48:DR:17:ARG:HE	1.83	0.43
48:DR:53:HIS:O	48:DR:56:LYS:CB	2.65	0.43
36:DB:7:G:C4'	49:DS:29:PHE:HE2	2.31	0.43
50:DT:107:ASP:CG	50:DT:109:GLU:H	2.20	0.43
51:DU:90:VAL:O	51:DU:92:ARG:N	2.48	0.43
52:DV:22:VAL:HB	52:DV:94:LEU:CB	2.39	0.43
52:DV:89:GLN:NE2	52:DV:89:GLN:HA	2.33	0.43
54:DX:88:LYS:HB3	54:DX:89:ILE:HD12	1.99	0.43
55:DY:28:LYS:HZ2	55:DY:37:VAL:CG1	2.26	0.43
55:DY:37:VAL:O	55:DY:38:ILE:CB	2.65	0.43
56:DZ:29:TYR:HA	56:DZ:33:LEU:O	2.18	0.43
56:DZ:37:VAL:O	56:DZ:37:VAL:HG23	2.17	0.43
1:AA:1151:A:C4	1:AA:1152:A:N7	2.86	0.43
1:AA:1387:G:C6	1:AA:1388:C:N4	2.87	0.43
1:AA:290:C:O2'	1:AA:291:C:H5'	2.18	0.43
1:AA:302:G:N3	1:AA:556:C:H4'	2.33	0.43
1:AA:452:A:H4'	16:AP:72:ARG:CZ	2.48	0.43
1:AA:50:A:N3	1:AA:52:G:H1'	2.34	0.43
1:AA:552:U:H2'	1:AA:553:A:C8	2.53	0.43
1:AA:608:A:H2'	1:AA:609:A:O4'	2.18	0.43
1:AA:640:A:H2'	1:AA:641:U:H5'	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:722:A:O2'	1:AA:723:U:C6	2.70	0.43
1:AA:780:A:H2	1:AA:803:G:C6	2.36	0.43
1:AA:969:A:O2'	1:AA:970:C:H5'	2.17	0.43
1:AA:973:G:C8	1:AA:974:A:C8	3.06	0.43
2:AB:22:LYS:O	2:AB:24:TRP:N	2.51	0.43
3:AC:172:ARG:NH1	3:AC:172:ARG:HB3	2.33	0.43
4:AD:105:VAL:HG13	4:AD:110:PHE:HB2	2.01	0.43
4:AD:150:GLU:O	4:AD:153:ARG:N	2.48	0.43
4:AD:157:LEU:O	4:AD:161:ASN:OD1	2.35	0.43
4:AD:161:ASN:O	4:AD:165:MET:HB2	2.18	0.43
4:AD:30:LYS:HA	4:AD:35:ARG:HD2	1.99	0.43
5:AE:150:ARG:CZ	5:AE:150:ARG:HB2	2.46	0.43
6:AF:72:VAL:O	6:AF:75:LEU:HB2	2.18	0.43
7:AG:25:ALA:CA	7:AG:28:ASN:HD22	2.31	0.43
8:AH:127:LEU:HD13	8:AH:127:LEU:C	2.39	0.43
8:AH:14:ARG:HH11	8:AH:14:ARG:HB3	1.82	0.43
8:AH:36:LEU:O	8:AH:37:ARG:C	2.57	0.43
8:AH:50:ARG:H	8:AH:50:ARG:HG3	1.61	0.43
11:AK:86:GLY:N	11:AK:112:THR:HG23	2.32	0.43
11:AK:92:GLU:HA	11:AK:95:ILE:HG12	2.01	0.43
12:AL:28:LYS:O	12:AL:30:ALA:N	2.51	0.43
13:AM:19:LEU:C	13:AM:22:ILE:HD13	2.39	0.43
14:AN:3:ARG:HB3	14:AN:3:ARG:NH1	2.32	0.43
15:AO:11:VAL:O	15:AO:14:GLU:HB3	2.18	0.43
16:AP:27:LYS:N	16:AP:27:LYS:HD2	2.28	0.43
19:AS:10:PHE:HE2	19:AS:70:LYS:NZ	2.14	0.43
1:AA:1328:C:P	21:AU:21:TYR:HH	2.42	0.43
1:AA:1340:A:OP2	23:AW:36:A:H5'	2.17	0.43
23:AW:39:A:H2'	23:AW:40:C:C5'	2.47	0.43
23:AW:49:C:H2'	23:AW:60:A:H4'	1.99	0.43
25:AY:132:ILE:O	25:AY:136:ALA:HB2	2.19	0.43
28:B2:20:GLU:O	28:B2:23:LYS:HB3	2.18	0.43
28:B2:57:ILE:CD1	28:B2:59:ARG:CZ	2.96	0.43
29:B3:14:GLY:H	29:B3:20:LYS:NZ	2.15	0.43
29:B3:9:VAL:HG23	29:B3:10:LYS:N	2.33	0.43
31:B5:15:ARG:HH11	31:B5:15:ARG:HG3	1.83	0.43
31:B5:15:ARG:HA	31:B5:18:ALA:HB2	1.98	0.43
32:B6:44:ARG:HG2	32:B6:44:ARG:HH11	1.84	0.43
34:B8:11:LYS:O	34:B8:11:LYS:CG	2.67	0.43
34:B8:6:THR:CB	34:B8:63:PRO:HD3	2.48	0.43
35:BA:1011:G:C5	35:BA:1151:G:N1	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1164:G:C6	35:BA:1165:U:C4	3.07	0.43
35:BA:121:G:H2'	35:BA:122:G:C8	2.52	0.43
35:BA:1309:G:H2'	35:BA:1310:G:C5'	2.48	0.43
35:BA:1469:A:C2'	35:BA:1470:G:H5'	2.47	0.43
35:BA:1668:A:C5	35:BA:1674:G:C5	3.06	0.43
35:BA:154:G:C6	35:BA:173:G:C6	3.06	0.43
35:BA:1766:U:H2'	35:BA:1767:C:C6	2.52	0.43
35:BA:2173:A:P	35:BA:2173:A:H3'	2.58	0.43
35:BA:2586:C:O2'	35:BA:2587:A:H5'	2.18	0.43
35:BA:1638:C:H1'	35:BA:2698:U:O2'	2.18	0.43
35:BA:410:G:C2	35:BA:418:G:C2	3.05	0.43
35:BA:431:U:C2'	35:BA:432:A:H5'	2.48	0.43
35:BA:606:U:H2'	35:BA:607:U:O4'	2.17	0.43
35:BA:723:G:C6	35:BA:724:U:C4	3.06	0.43
35:BA:869:G:C4	35:BA:870:A:C8	3.06	0.43
35:BA:975(A):G:N3	35:BA:1156:A:H2	2.15	0.43
37:BC:21:THR:O	37:BC:22:ILE:C	2.57	0.43
38:BD:133:LEU:HB3	38:BD:173:VAL:HG11	2.00	0.43
39:BE:188:VAL:O	39:BE:188:VAL:HG13	2.18	0.43
39:BE:49:LEU:HD23	39:BE:81:ILE:CG1	2.48	0.43
39:BE:53:PRO:O	39:BE:55:ASN:OD1	2.36	0.43
41:BG:27:ASN:CG	41:BG:28:VAL:N	2.72	0.43
42:BH:83:TYR:HD1	42:BH:84:SER:N	2.16	0.43
44:BN:110:GLY:O	44:BN:113:GLY:N	2.51	0.43
44:BN:23:LEU:O	44:BN:25:ARG:N	2.48	0.43
44:BN:26:LEU:CD2	44:BN:99:LEU:HD11	2.48	0.43
44:BN:36:GLY:HA3	44:BN:48:MET:SD	2.58	0.43
44:BN:39:ARG:HG3	44:BN:39:ARG:HH11	1.83	0.43
44:BN:42:TRP:HE3	44:BN:48:MET:SD	2.41	0.43
44:BN:66:LYS:HA	44:BN:66:LYS:HE3	2.01	0.43
47:BQ:127:ILE:HG22	47:BQ:128:LYS:N	2.19	0.43
48:BR:4:LEU:HD22	48:BR:4:LEU:O	2.18	0.43
48:BR:4:LEU:O	48:BR:6:SER:N	2.51	0.43
35:BA:2880:C:H1'	48:BR:92:GLY:O	2.18	0.43
49:BS:66:ALA:CA	49:BS:69:VAL:HG12	2.48	0.43
51:BU:59:ARG:C	51:BU:61:TRP:N	2.68	0.43
51:BU:88:ILE:O	51:BU:88:ILE:CD1	2.66	0.43
53:BW:84:ARG:HB2	53:BW:96:ILE:CG2	2.48	0.43
54:BX:35:THR:O	54:BX:36:LYS:O	2.36	0.43
55:BY:84:ARG:C	55:BY:85:VAL:CG2	2.85	0.43
56:BZ:167:PRO:O	56:BZ:168:GLU:CB	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1164:G:C2'	1:CA:1165:C:H5'	2.49	0.43
1:CA:1220:G:H21	19:CS:54:GLY:HA2	1.83	0.43
1:CA:189(B):C:H2'	1:CA:189(C):C:H6	1.82	0.43
1:CA:355:C:C2	1:CA:356:A:C8	3.06	0.43
1:CA:381:C:H2'	1:CA:382:A:O4'	2.18	0.43
1:CA:460:G:N2	1:CA:472:A:H62	2.15	0.43
1:CA:521:G:O2'	1:CA:522:C:H5'	2.18	0.43
1:CA:544:G:C5	1:CA:545:C:C5	3.06	0.43
1:CA:61:G:O2'	1:CA:62:U:H5'	2.18	0.43
1:CA:763:G:C5	1:CA:764:C:C5	3.06	0.43
1:CA:866:C:C2	1:CA:867:G:H1'	2.53	0.43
1:CA:953:G:O2'	1:CA:954:G:H5'	2.18	0.43
2:CB:115:LEU:HD21	2:CB:153:ARG:NE	2.32	0.43
2:CB:212:GLN:NE2	2:CB:216:SER:CB	2.78	0.43
2:CB:36:ARG:HG3	2:CB:37:ASN:N	2.33	0.43
3:CC:150:LYS:O	3:CC:201:TYR:N	2.46	0.43
3:CC:25:GLY:O	3:CC:28:GLN:N	2.51	0.43
4:CD:13:ARG:O	4:CD:14:ARG:C	2.56	0.43
4:CD:58:LEU:C	4:CD:58:LEU:HD13	2.38	0.43
7:CG:62:PHE:CG	7:CG:62:PHE:O	2.71	0.43
8:CH:120:THR:C	8:CH:122:ARG:H	2.21	0.43
8:CH:12:ARG:HA	8:CH:15:ASN:HD22	1.82	0.43
8:CH:51:VAL:HG11	8:CH:60:ARG:CG	2.45	0.43
8:CH:85:ARG:HH11	8:CH:85:ARG:HG3	1.83	0.43
9:CI:116:LYS:HA	9:CI:121:ARG:O	2.18	0.43
9:CI:125:TYR:C	9:CI:125:TYR:CD2	2.91	0.43
10:CJ:3:LYS:O	10:CJ:100:THR:HA	2.19	0.43
10:CJ:23:ILE:HG23	10:CJ:85:LEU:HD22	2.00	0.43
11:CK:65:ALA:HB1	11:CK:98:LEU:HD23	2.01	0.43
1:CA:994:A:H2	14:CN:4:LYS:HG3	1.83	0.43
15:CO:3:ILE:HG13	15:CO:3:ILE:O	2.18	0.43
15:CO:70:LEU:HD12	15:CO:70:LEU:HA	1.85	0.43
16:CP:77:ALA:HB3	16:CP:79:VAL:HG23	2.00	0.43
17:CQ:11:VAL:HG23	17:CQ:20:THR:CG2	2.49	0.43
17:CQ:59:ILE:HG23	17:CQ:71:PHE:HB3	2.00	0.43
17:CQ:74:LEU:HD12	17:CQ:75:ARG:CG	2.41	0.43
23:CW:25:U:H2'	23:CW:26:C:O4'	2.18	0.43
25:CY:105:PRO:O	25:CY:106:LEU:HD23	2.19	0.43
25:CY:171:LYS:HD3	25:CY:171:LYS:HA	1.83	0.43
27:D1:85:LEU:CD2	27:D1:85:LEU:H	2.31	0.43
28:D2:15:LYS:O	28:D2:16:LEU:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D2:41:ILE:O	28:D2:43:GLN:N	2.51	0.43
29:D3:35:ARG:HG3	29:D3:35:ARG:NH1	2.33	0.43
33:D7:34:ARG:O	33:D7:36:GLN:N	2.52	0.43
34:D8:29:LYS:NZ	34:D8:44:LYS:HB2	2.33	0.43
34:D8:56:GLU:O	34:D8:59:LYS:NZ	2.39	0.43
35:DA:103:A:H2'	35:DA:104:U:C6	2.53	0.43
35:DA:1142(A):A:C4	35:DA:1144:G:N7	2.86	0.43
35:DA:1187:G:OP1	52:DV:82:ARG:NH2	2.40	0.43
35:DA:1263:U:C5	35:DA:1264:G:C6	3.06	0.43
35:DA:1281:G:H1	35:DA:1286:A:H62	1.65	0.43
35:DA:1525:G:H2'	35:DA:1526:G:C8	2.53	0.43
35:DA:1823:G:O2'	35:DA:1824:G:H5'	2.18	0.43
35:DA:1916:A:H3'	35:DA:1917:U:H6	1.82	0.43
35:DA:2008:C:H2'	35:DA:2009:G:H8	1.82	0.43
35:DA:2199:A:N3	35:DA:2199:A:H2'	2.33	0.43
35:DA:260:G:N3	35:DA:260:G:H2'	2.32	0.43
35:DA:2821:A:H2'	35:DA:2822:G:H8	1.76	0.43
35:DA:510:C:OP1	35:DA:511:U:OP2	2.36	0.43
35:DA:564:C:H2'	35:DA:565:C:H6	1.75	0.43
35:DA:605:C:H2'	35:DA:606:U:H6	1.83	0.43
35:DA:619:G:H5''	35:DA:620:G:OP2	2.18	0.43
35:DA:598:G:C6	35:DA:660:G:C6	3.06	0.43
35:DA:70:G:H21	35:DA:71:A:H62	1.65	0.43
35:DA:724:U:C2'	35:DA:725:G:H5'	2.48	0.43
35:DA:74:A:O2'	35:DA:75:G:OP2	2.32	0.43
35:DA:804:A:H2'	35:DA:806:C:N4	2.33	0.43
38:DD:134:ARG:C	38:DD:136:ILE:H	2.21	0.43
38:DD:34:VAL:HG22	38:DD:35:LYS:NZ	2.33	0.43
39:DE:120:TRP:CD2	39:DE:155:LYS:HD3	2.53	0.43
39:DE:132:HIS:O	39:DE:135:HIS:CD2	2.72	0.43
35:DA:2051:A:C4'	39:DE:141:ILE:HD11	2.47	0.43
39:DE:3:GLY:O	39:DE:4:ILE:CB	2.66	0.43
39:DE:9:VAL:CG2	50:DT:8:LYS:HB2	2.49	0.43
40:DF:155:LEU:HD12	40:DF:174:VAL:O	2.18	0.43
40:DF:9:ILE:HG12	40:DF:15:SER:N	2.32	0.43
35:DA:323:G:C3'	40:DF:169:ASN:HD21	2.30	0.43
40:DF:6:VAL:HB	40:DF:7:TYR:H	1.62	0.43
35:DA:2306:C:N4	41:DG:43:LEU:O	2.43	0.43
43:DI:54:GLN:HA	43:DI:57:ARG:HB3	2.00	0.43
44:DN:62:VAL:HG21	44:DN:66:LYS:HB2	2.00	0.43
47:DQ:77:LYS:HA	47:DQ:78:PRO:HD3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:DR:62:ALA:O	48:DR:63:ARG:C	2.56	0.43
51:DU:14:HIS:C	51:DU:16:LYS:N	2.70	0.43
52:DV:3:ALA:O	52:DV:14:VAL:N	2.46	0.43
53:DW:14:PRO:O	53:DW:15:ARG:C	2.56	0.43
54:DX:36:LYS:HZ2	54:DX:39:ILE:HA	1.78	0.43
54:DX:43:VAL:C	54:DX:45:THR:N	2.71	0.43
55:DY:34:LYS:O	55:DY:35:TYR:HB3	2.18	0.43
55:DY:14:LEU:O	55:DY:72:VAL:HA	2.17	0.43
56:DZ:30:ASN:O	56:DZ:32:HIS:N	2.51	0.43
1:AA:1169:A:C2	1:AA:1170:A:C4	3.06	0.43
1:AA:1365:G:H2'	1:AA:1366:C:O4'	2.17	0.43
1:AA:1489:G:C4	1:AA:1490:C:C5	3.06	0.43
1:AA:189(E):U:O2'	1:AA:189(F):U:H5'	2.18	0.43
1:AA:23:C:C2'	1:AA:24:U:H5'	2.48	0.43
1:AA:267:C:OP2	17:AQ:67:LYS:HD2	2.18	0.43
1:AA:44:G:C2	1:AA:399:G:C2	3.07	0.43
1:AA:544:G:C5	1:AA:545:C:C5	3.06	0.43
1:AA:582:U:H2'	1:AA:583:A:H8	1.83	0.43
1:AA:779:C:O2'	11:AK:120:ARG:HD3	2.18	0.43
2:AB:180:LEU:C	2:AB:182:ILE:H	2.21	0.43
4:AD:177:ASP:O	4:AD:177:ASP:OD1	2.36	0.43
4:AD:92:VAL:O	4:AD:93:PHE:C	2.56	0.43
1:AA:933:G:OP2	7:AG:3:ARG:HB3	2.19	0.43
8:AH:53:VAL:HG12	8:AH:54:ASP:OD2	2.19	0.43
9:AI:17:VAL:HA	9:AI:63:ILE:HG23	1.99	0.43
12:AL:27:LEU:C	12:AL:29:GLY:N	2.71	0.43
12:AL:95:GLY:O	12:AL:97:ARG:N	2.51	0.43
13:AM:124:PRO:HB2	13:AM:125:ARG:H	1.63	0.43
10:AJ:49:VAL:HG11	14:AN:41:ARG:HB2	2.01	0.43
18:AR:62:GLU:O	18:AR:63:GLN:C	2.57	0.43
19:AS:36:ARG:HH11	19:AS:36:ARG:HB3	1.83	0.43
20:AT:97:ALA:HA	20:AT:98:PRO:HD3	1.87	0.43
25:AY:29:ARG:HA	25:AY:32:ARG:HE	1.83	0.43
25:AY:87:ASP:O	25:AY:88:LEU:C	2.56	0.43
26:B0:75:LEU:HD23	26:B0:75:LEU:O	2.18	0.43
25:AY:150:SER:OG	26:B0:7:LEU:CB	2.67	0.43
28:B2:51:ARG:HD3	28:B2:51:ARG:C	2.39	0.43
29:B3:19:GLN:C	29:B3:21:ALA:H	2.21	0.43
29:B3:22:ALA:O	29:B3:26:LEU:HG	2.18	0.43
29:B3:6:VAL:HG12	29:B3:56:VAL:HA	2.00	0.43
34:B8:59:LYS:CB	34:B8:59:LYS:NZ	2.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1022:G:O2'	35:BA:1023:U:OP2	2.32	0.43
35:BA:1035:U:H5''	42:BH:59:ARG:NH1	2.33	0.43
35:BA:1195:G:C2	35:BA:1196:C:C5	3.07	0.43
35:BA:1526:G:H2'	35:BA:1527:G:O4'	2.18	0.43
35:BA:2487:G:O2'	35:BA:2488:A:H5'	2.18	0.43
35:BA:2495:G:C6	35:BA:2496:C:C4	3.06	0.43
35:BA:2714:G:C6	35:BA:2715:C:N3	2.87	0.43
35:BA:2727:G:C5	35:BA:2728:U:H5	2.36	0.43
35:BA:2755:C:H6	35:BA:2755:C:H3'	1.83	0.43
35:BA:322:A:OP1	40:BF:168:ARG:HD3	2.19	0.43
35:BA:272(C):G:H1	35:BA:365:C:H42	1.66	0.43
33:B7:34:ARG:NH1	35:BA:466:A:OP1	2.52	0.43
35:BA:562:U:C2'	35:BA:563:G:OP2	2.66	0.43
35:BA:708:C:H5'	35:BA:709:U:OP2	2.18	0.43
36:BB:70:C:O2'	36:BB:71:C:H5'	2.19	0.43
36:BB:74:U:C5	36:BB:75:G:N7	2.87	0.43
36:BB:81:G:H4'	36:BB:81:G:OP1	2.18	0.43
37:BC:184:LYS:C	37:BC:186:ALA:N	2.71	0.43
38:BD:53:PHE:HB3	38:BD:218:ARG:HB2	2.00	0.43
39:BE:103:ASP:OD1	39:BE:201:THR:HA	2.19	0.43
39:BE:79:ARG:NH1	39:BE:79:ARG:HG2	2.32	0.43
40:BF:26:ALA:HB1	40:BF:27:GLU:OE1	2.18	0.43
40:BF:32:LEU:O	40:BF:35:GLU:CB	2.67	0.43
40:BF:45:ARG:NH1	40:BF:97:TYR:CE2	2.87	0.43
41:BG:174:GLU:C	41:BG:176:LEU:H	2.22	0.43
41:BG:69:ALA:O	41:BG:90:LEU:HA	2.18	0.43
43:BI:93:THR:N	43:BI:96:ASP:OD2	2.51	0.43
44:BN:46:VAL:CG2	44:BN:48:MET:HG3	2.48	0.43
45:BO:1:MET:H3	45:BO:1:MET:CE	2.30	0.43
46:BP:107:LYS:C	46:BP:109:GLY:N	2.71	0.43
46:BP:95:VAL:HG23	46:BP:125:VAL:CG2	2.48	0.43
47:BQ:81:VAL:CG2	47:BQ:82:ARG:HG2	2.47	0.43
48:BR:20:LEU:O	48:BR:21:TYR:C	2.54	0.43
48:BR:34:ILE:HG22	48:BR:35:THR:H	1.84	0.43
49:BS:106:ARG:NH1	49:BS:107:GLU:O	2.51	0.43
49:BS:26:LEU:N	49:BS:88:ASP:HB2	2.33	0.43
50:BT:35:LYS:NZ	50:BT:41:ARG:HE	2.13	0.43
50:BT:72:VAL:CG1	50:BT:73:GLU:N	2.81	0.43
51:BU:36:ARG:HG2	51:BU:40:PHE:CE2	2.52	0.43
53:BW:77:ASP:O	53:BW:102:HIS:HB2	2.18	0.43
55:BY:7:VAL:HB	55:BY:8:LYS:NZ	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1332:A:C2	1:CA:1333:A:C5	3.07	0.43
1:CA:1480:G:H2'	1:CA:1481:U:H6	1.83	0.43
1:CA:1497:G:HO2'	1:CA:1518:A:H2	1.65	0.43
1:CA:1518:A:N1	1:CA:1519:A:C6	2.86	0.43
1:CA:24:U:O2'	1:CA:25:C:H5'	2.19	0.43
1:CA:391:G:C6	1:CA:392:G:N7	2.86	0.43
1:CA:404:U:C2	1:CA:405:U:C5	3.07	0.43
1:CA:723:U:H5''	1:CA:724:G:OP2	2.18	0.43
1:CA:811:C:H4'	1:CA:900:A:H62	1.80	0.43
1:CA:908:A:C2'	1:CA:909:A:H8	2.08	0.43
2:CB:71:VAL:O	2:CB:164:VAL:HG13	2.17	0.43
3:CC:154:SER:O	3:CC:165:THR:HA	2.19	0.43
3:CC:92:ALA:HB2	3:CC:99:VAL:HG22	1.99	0.43
4:CD:79:PHE:HE1	4:CD:204:ILE:HA	1.82	0.43
4:CD:72:GLU:O	4:CD:76:ARG:HB2	2.19	0.43
5:CE:12:LEU:C	5:CE:12:LEU:HD22	2.39	0.43
6:CF:33:TYR:CE1	6:CF:75:LEU:HA	2.53	0.43
6:CF:54:LYS:O	6:CF:56:PRO:HD3	2.19	0.43
8:CH:111:ILE:N	8:CH:135:CYS:O	2.51	0.43
8:CH:28:ALA:O	8:CH:29:SER:HB2	2.18	0.43
11:CK:88:GLY:C	11:CK:91:ARG:HB2	2.38	0.43
18:CR:53:ARG:NH2	18:CR:60:ALA:CA	2.81	0.43
1:CA:1319:A:OP2	19:CS:5:LEU:HD23	2.18	0.43
20:CT:13:LEU:HD13	20:CT:14:LYS:N	2.33	0.43
20:CT:24:LEU:O	20:CT:27:LYS:N	2.52	0.43
1:CA:1397:C:N3	24:CX:22:U:C6	2.86	0.43
25:CY:123:GLU:O	25:CY:126:ARG:HB2	2.18	0.43
25:CY:152:ASP:O	25:CY:155:LYS:HB2	2.18	0.43
25:CY:45:TYR:HB2	25:CY:78:ALA:CB	2.46	0.43
26:D0:25:ARG:HG3	26:D0:29:GLN:HE21	1.81	0.43
28:D2:14:ARG:HE	28:D2:15:LYS:N	2.15	0.43
32:D6:40:CYS:SG	32:D6:45:LYS:HD2	2.59	0.43
34:D8:32:LEU:HD23	34:D8:35:GLN:C	2.39	0.43
34:D8:35:GLN:HE21	34:D8:36:LYS:HZ3	1.66	0.43
35:DA:1174:A:OP1	35:DA:1175:U:H5''	2.19	0.43
35:DA:1306:C:H2'	35:DA:1307:A:H8	1.83	0.43
35:DA:1474:C:H3'	35:DA:1475:G:H8	1.83	0.43
35:DA:1803:A:H4'	38:DD:259:THR:CG2	2.48	0.43
35:DA:191:A:C2	35:DA:192:C:C4	3.07	0.43
35:DA:2020:A:C5	35:DA:2022:U:C5	3.06	0.43
35:DA:2300:G:H1	35:DA:2316:C:N4	2.10	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:243:U:C2'	35:DA:244:A:H5'	2.48	0.43
35:DA:2566:A:C6	45:DO:28:SER:HB2	2.53	0.43
35:DA:2574:G:C5	35:DA:2575:C:C5	3.07	0.43
35:DA:2886:G:H2'	35:DA:2887:U:C6	2.54	0.43
35:DA:491:G:C4	35:DA:492:A:C8	3.06	0.43
35:DA:514:A:C2	35:DA:515:A:C4	3.06	0.43
35:DA:769:G:H2'	35:DA:770:G:H8	1.83	0.43
35:DA:869:G:C4	35:DA:870:A:C8	3.06	0.43
35:DA:884:C:H4'	35:DA:892:G:C8	2.54	0.43
35:DA:994:C:H1'	52:DV:10:LYS:HE2	2.00	0.43
36:DB:81:G:H5'	36:DB:82:G:OP2	2.18	0.43
38:DD:145:VAL:HG12	38:DD:146:GLU:H	1.82	0.43
38:DD:199:ALA:O	38:DD:201:HIS:N	2.52	0.43
38:DD:75:ILE:HG21	38:DD:99:ASP:HB2	2.01	0.43
39:DE:49:LEU:N	39:DE:49:LEU:CD2	2.77	0.43
41:DG:136:ARG:O	41:DG:136:ARG:HD3	2.17	0.43
41:DG:14:GLU:OE1	41:DG:15:VAL:HG23	2.17	0.43
35:DA:2313:C:OP1	41:DG:71:THR:HG21	2.19	0.43
41:DG:96:ARG:HA	41:DG:99:MET:HE2	2.01	0.43
42:DH:86:GLU:HA	42:DH:132:ARG:CB	2.48	0.43
43:DI:109:ILE:HD12	43:DI:111:PRO:HD3	2.00	0.43
45:DO:113:LYS:HA	45:DO:116:SER:OG	2.19	0.43
45:DO:1:MET:CG	45:DO:32:TYR:CD2	2.95	0.43
35:DA:1245:G:OP1	46:DP:16:ARG:NE	2.51	0.43
47:DQ:48:GLU:O	47:DQ:49:ALA:C	2.56	0.43
48:DR:49:ASP:O	48:DR:52:ILE:N	2.51	0.43
48:DR:60:LEU:HA	48:DR:63:ARG:HB2	2.00	0.43
48:DR:73:VAL:HG23	48:DR:74:LYS:N	2.33	0.43
49:DS:26:LEU:N	49:DS:88:ASP:HB2	2.34	0.43
50:DT:51:ARG:O	50:DT:61:PHE:HA	2.19	0.43
52:DV:34:GLU:O	52:DV:62:LEU:HG	2.19	0.43
52:DV:75:PHE:HD1	52:DV:87:HIS:HB3	1.81	0.43
35:DA:493:G:HO2'	53:DW:8:ARG:H	1.65	0.43
54:DX:53:LYS:CE	54:DX:55:ASN:HD21	2.32	0.43
54:DX:77:LYS:HA	54:DX:77:LYS:CE	2.46	0.43
56:DZ:116:VAL:HG12	56:DZ:117:LEU:N	2.31	0.43
56:DZ:5:LEU:O	56:DZ:59:LEU:HA	2.19	0.43
56:DZ:6:LYS:CB	56:DZ:8:TYR:CE1	2.98	0.43
1:AA:1065:U:O2'	1:AA:1066:C:OP2	2.26	0.43
1:AA:107:G:H2'	1:AA:108:G:H5'	2.00	0.43
1:AA:1296:C:H4'	1:AA:1302:U:O4	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1423:G:C6	1:AA:1424:C:C4	3.06	0.43
1:AA:152:A:N6	1:AA:170:U:C2	2.87	0.43
1:AA:246:A:O3'	1:AA:247:G:H4'	2.19	0.43
1:AA:741:G:C2'	1:AA:742:G:O4'	2.66	0.43
1:AA:956:U:H2'	1:AA:957:U:C6	2.54	0.43
2:AB:171:ALA:O	2:AB:174:VAL:N	2.51	0.43
3:AC:173:VAL:O	3:AC:173:VAL:CG1	2.63	0.43
4:AD:148:VAL:O	4:AD:149:ALA:O	2.36	0.43
4:AD:39:PRO:HB3	4:AD:40:PRO:HD2	2.01	0.43
5:AE:41:VAL:CG1	5:AE:113:ALA:HA	2.49	0.43
5:AE:6:PHE:HB2	5:AE:34:VAL:CG1	2.48	0.43
7:AG:111:ARG:HB3	7:AG:111:ARG:HH11	1.82	0.43
7:AG:60:LYS:NZ	7:AG:63:LYS:HG2	2.34	0.43
7:AG:92:SER:CB	7:AG:94:ARG:HE	2.31	0.43
9:AI:86:VAL:HG13	9:AI:87:GLN:N	2.33	0.43
1:AA:692:U:H5	11:AK:26:ASN:CG	2.22	0.43
25:AY:150:SER:HB2	25:AY:152:ASP:OD1	2.18	0.43
25:AY:174:GLN:NE2	25:AY:175:LEU:N	2.67	0.43
25:AY:64:ARG:NE	25:AY:64:ARG:HA	2.34	0.43
26:B0:32:ARG:HB2	26:B0:35:ASN:ND2	2.34	0.43
32:B6:15:GLU:CD	32:B6:18:ARG:HG3	2.39	0.43
35:BA:1406:U:H2'	35:BA:1407:C:C6	2.54	0.43
35:BA:1511:C:H2'	35:BA:1512:U:C6	2.53	0.43
35:BA:1586:A:N1	35:BA:1587:A:C4	2.86	0.43
35:BA:1632:A:C6	35:BA:1633:G:C6	3.06	0.43
35:BA:49:A:C6	35:BA:177:G:C5	3.06	0.43
35:BA:1800:C:N3	35:BA:1818:U:O2	2.51	0.43
35:BA:2205:C:C2	35:BA:2220:G:N1	2.87	0.43
35:BA:225:A:H2'	35:BA:226:G:O4'	2.18	0.43
35:BA:2283:C:C6	35:BA:2389:G:H2'	2.54	0.43
35:BA:2307:G:C3'	35:BA:2307:G:N3	2.78	0.43
35:BA:258:G:C6	35:BA:259:G:C5	3.06	0.43
35:BA:2801:A:O2'	35:BA:2895:U:C4'	2.66	0.43
35:BA:486:C:O2	35:BA:495:G:C2	2.72	0.43
35:BA:880:G:N1	35:BA:898:C:N4	2.66	0.43
35:BA:998:C:OP2	51:BU:93:LYS:NZ	2.48	0.43
36:BB:24:G:H4'	36:BB:25:A:N7	2.34	0.43
36:BB:43:C:H3'	36:BB:44:G:H5'	2.00	0.43
37:BC:82:LYS:O	37:BC:83:ILE:HD13	2.19	0.43
38:BD:76:PRO:O	38:BD:98:VAL:HG23	2.18	0.43
39:BE:14:ILE:O	39:BE:20:ALA:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:10:PRO:CG	40:BF:13:SER:OG	2.67	0.43
41:BG:138:GLN:HG3	41:BG:152:LEU:HD22	2.00	0.43
41:BG:5:VAL:CG1	41:BG:6:ALA:H	2.08	0.43
42:BH:149:ARG:HH21	42:BH:154:PRO:HG3	1.84	0.43
43:BI:29:TYR:C	43:BI:32:PRO:HD2	2.39	0.43
44:BN:15:LEU:C	44:BN:15:LEU:HD13	2.38	0.43
45:BO:104:ARG:HH21	50:BT:33:LYS:CE	2.31	0.43
45:BO:64:ARG:HB2	45:BO:64:ARG:HH11	1.84	0.43
47:BQ:90:VAL:HG12	47:BQ:90:VAL:O	2.18	0.43
49:BS:37:ALA:O	49:BS:38:GLN:HB2	2.18	0.43
49:BS:69:VAL:CG1	49:BS:70:GLY:N	2.81	0.43
51:BU:8:VAL:O	51:BU:9:VAL:C	2.57	0.43
52:BV:1:MET:CE	52:BV:46:VAL:HG23	2.47	0.43
53:BW:27:LYS:H	53:BW:71:VAL:HB	1.82	0.43
54:BX:60:ARG:CB	54:BX:72:LYS:H	2.30	0.43
56:BZ:140:ASP:O	56:BZ:142:SER:N	2.51	0.43
1:CA:1095:U:H2'	1:CA:1096:C:H6	1.82	0.43
1:CA:1089:G:H1	1:CA:1096:C:H42	1.66	0.43
1:CA:1105:A:O2'	1:CA:1106:G:H5'	2.18	0.43
1:CA:1296:C:H3'	1:CA:1297:C:C6	2.52	0.43
1:CA:1402:C:H2'	1:CA:1403:C:C5'	2.49	0.43
1:CA:27:G:H2'	1:CA:28:G:O4'	2.18	0.43
1:CA:376:G:O3'	16:CP:5:ARG:HD2	2.18	0.43
1:CA:407:G:N1	1:CA:408:A:C5	2.87	0.43
1:CA:445:G:H2'	1:CA:446:G:O4'	2.19	0.43
1:CA:44:G:H2'	1:CA:45:U:O4'	2.18	0.43
1:CA:584:G:O2'	1:CA:585:G:H5'	2.19	0.43
1:CA:731:G:O2'	1:CA:732:C:H5'	2.18	0.43
1:CA:890:G:N2	1:CA:906:G:H2'	2.33	0.43
2:CB:77:ALA:HB1	2:CB:211:ILE:HD13	2.00	0.43
7:CG:18:TYR:HE1	7:CG:44:TYR:CE2	2.36	0.43
9:CI:8:GLY:O	9:CI:14:VAL:HG13	2.18	0.43
10:CJ:22:LYS:C	10:CJ:22:LYS:HD2	2.39	0.43
1:CA:1253:G:OP1	10:CJ:44:VAL:HG11	2.18	0.43
11:CK:30:VAL:HG23	11:CK:30:VAL:O	2.18	0.43
11:CK:79:SER:HA	11:CK:104:GLN:HB3	2.00	0.43
11:CK:92:GLU:HA	11:CK:95:ILE:HG12	2.00	0.43
13:CM:36:LYS:O	13:CM:37:THR:HG23	2.17	0.43
14:CN:34:TYR:O	14:CN:36:PHE:N	2.52	0.43
15:CO:32:LEU:O	15:CO:36:ILE:HG13	2.18	0.43
17:CQ:60:ILE:HB	17:CQ:74:LEU:HD23	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1320:C:H5'	19:CS:70:LYS:HD3	2.00	0.43
21:CU:7:ARG:O	21:CU:21:TYR:HD2	2.01	0.43
23:CW:2:G:H21	23:CW:3:C:H1'	1.80	0.43
23:CW:8:U:O4	23:CW:14:A:N7	2.51	0.43
25:CY:180:GLU:O	25:CY:181:GLN:C	2.56	0.43
25:CY:3:LEU:O	25:CY:5:GLU:N	2.51	0.43
28:D2:47:ASN:O	28:D2:48:HIS:C	2.56	0.43
34:D8:29:LYS:HZ1	34:D8:44:LYS:HB3	1.84	0.43
34:D8:22:VAL:CB	34:D8:53:PRO:HB2	2.36	0.43
34:D8:4:MET:CE	34:D8:61:LEU:HD12	2.48	0.43
35:DA:1128:A:H2	35:DA:2516:G:N3	2.15	0.43
35:DA:1289:C:C2	35:DA:1290:C:C5	3.07	0.43
35:DA:1340:U:C5	35:DA:1603:A:C1'	3.01	0.43
35:DA:1507:A:C2	35:DA:1508:A:H1'	2.54	0.43
35:DA:1835:G:H2'	35:DA:1835:G:N3	2.33	0.43
35:DA:1861:G:H2'	35:DA:1862:G:H8	1.83	0.43
35:DA:186:G:C2'	35:DA:187:G:H5'	2.48	0.43
35:DA:1916:A:H3'	35:DA:1917:U:C6	2.53	0.43
35:DA:207:A:H2'	35:DA:208:C:O4'	2.19	0.43
35:DA:2228:G:H2'	35:DA:2229:C:C6	2.53	0.43
35:DA:2322:A:H2'	35:DA:2323:G:O4'	2.18	0.43
35:DA:2483:C:O2	35:DA:2483:C:H2'	2.18	0.43
35:DA:2488:A:O2'	35:DA:2489:G:H5'	2.18	0.43
35:DA:2718:G:O2'	35:DA:2719:G:H5'	2.18	0.43
35:DA:2633:G:H5'	35:DA:2811:G:O2'	2.18	0.43
35:DA:2824:C:H2'	35:DA:2825:C:O4'	2.19	0.43
35:DA:404:C:C3'	35:DA:405:U:H5'	2.47	0.43
35:DA:738:G:C2'	35:DA:739:G:H5'	2.48	0.43
35:DA:785:G:C5	35:DA:786:C:C5	3.06	0.43
35:DA:827:U:H2'	35:DA:2068:U:H3	1.83	0.43
35:DA:93:G:H2'	35:DA:94:C:O4'	2.18	0.43
36:DB:99:G:O2'	36:DB:100:A:O4'	2.30	0.43
38:DD:139:GLY:O	38:DD:140:THR:C	2.57	0.43
38:DD:142:VAL:HA	38:DD:194:GLY:N	2.32	0.43
38:DD:235:GLY:O	38:DD:237:GLU:HG2	2.19	0.43
38:DD:271:ILE:O	38:DD:272:ALA:O	2.37	0.43
38:DD:57:GLY:O	38:DD:58:HIS:O	2.35	0.43
39:DE:108:SER:O	39:DE:162:ALA:HA	2.17	0.43
40:DF:170:LEU:HD23	40:DF:173:VAL:HB	2.00	0.43
40:DF:22:ALA:HB1	40:DF:26:ALA:CB	2.47	0.43
40:DF:2:LYS:HG3	40:DF:25:PRO:CB	2.23	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DF:63:LYS:HA	40:DF:76:GLY:O	2.19	0.43
41:DG:139:LEU:HD22	41:DG:146:TYR:CD1	2.53	0.43
41:DG:161:THR:HG22	41:DG:163:ALA:CB	2.49	0.43
42:DH:127:GLU:CB	42:DH:130:ARG:HB3	2.49	0.43
42:DH:160:LYS:O	42:DH:163:TYR:CE1	2.71	0.43
42:DH:89:ILE:HD13	42:DH:90:LYS:N	2.33	0.43
44:DN:55:VAL:CG1	44:DN:126:PRO:HA	2.48	0.43
45:DO:115:VAL:O	45:DO:118:ALA:HB3	2.19	0.43
45:DO:61:VAL:O	45:DO:61:VAL:HG13	2.18	0.43
34:D8:30:ARG:HH21	46:DP:62:LEU:HD23	1.82	0.43
46:DP:96:THR:HB	46:DP:97:PRO:HD2	2.01	0.43
47:DQ:134:ARG:CG	47:DQ:135:ASP:H	2.15	0.43
47:DQ:16:ARG:CB	47:DQ:16:ARG:HH11	2.30	0.43
47:DQ:42:ILE:N	47:DQ:42:ILE:HD12	2.33	0.43
48:DR:99:LYS:C	48:DR:100:LEU:HD22	2.39	0.43
48:DR:31:HIS:O	48:DR:33:ARG:N	2.49	0.43
51:DU:92:ARG:HG3	51:DU:94:ASN:H	1.82	0.43
51:DU:92:ARG:CZ	52:DV:11:GLN:H	2.31	0.43
52:DV:70:ILE:CG2	52:DV:90:PRO:HB2	2.48	0.43
54:DX:75:ASP:O	54:DX:76:ARG:HG3	2.19	0.43
55:DY:86:ARG:NH2	55:DY:95:LYS:NZ	2.66	0.43
56:DZ:10:ARG:CG	56:DZ:38:TYR:HD2	2.30	0.43
56:DZ:45:ASP:C	56:DZ:49:ARG:HG2	2.38	0.43
1:AA:1098:C:C2	1:AA:1099:G:C8	3.06	0.43
1:AA:1146:A:C2'	1:AA:1147:C:O5'	2.67	0.43
1:AA:1161:C:H2'	1:AA:1162:C:C6	2.54	0.43
1:AA:1288:A:O5'	1:AA:1288:A:H8	2.01	0.43
1:AA:1368:G:H4'	14:AN:61:TRP:HZ2	1.83	0.43
1:AA:1476:G:C6	1:AA:1477:C:C4	3.07	0.43
1:AA:235:C:O2'	1:AA:236:G:H5'	2.19	0.43
1:AA:514:C:N4	1:AA:537:G:H1	2.15	0.43
1:AA:777:A:O2'	1:AA:778:G:H5'	2.19	0.43
1:AA:792:A:N3	1:AA:794:A:C5	2.87	0.43
1:AA:953:G:O2'	1:AA:954:G:H5'	2.19	0.43
1:AA:954:G:C5	1:AA:955:U:C4	3.07	0.43
2:AB:177:ALA:O	2:AB:180:LEU:N	2.49	0.43
3:AC:6:HIS:CD2	3:AC:8:ILE:H	2.36	0.43
3:AC:95:THR:HG22	3:AC:97:LYS:H	1.84	0.43
4:AD:194:LEU:O	4:AD:195:ALA:HB3	2.18	0.43
5:AE:72:GLN:OE1	5:AE:77:PRO:HA	2.18	0.43
6:AF:100:ASN:HD22	6:AF:100:ASN:HA	1.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:23:VAL:C	7:AG:27:ILE:HD13	2.38	0.43
7:AG:60:LYS:HD2	7:AG:63:LYS:HG2	1.99	0.43
8:AH:120:THR:N	8:AH:123:GLU:OE1	2.52	0.43
8:AH:26:VAL:HG12	8:AH:59:LEU:HB2	2.01	0.43
8:AH:46:LYS:HG3	8:AH:63:LEU:O	2.19	0.43
12:AL:102:ARG:NH1	12:AL:102:ARG:CG	2.78	0.43
13:AM:51:ALA:O	13:AM:55:ARG:HB2	2.18	0.43
14:AN:9:LYS:HB2	14:AN:9:LYS:HE3	1.82	0.43
15:AO:66:LEU:O	15:AO:69:TYR:N	2.51	0.43
17:AQ:74:LEU:HD12	17:AQ:75:ARG:CG	2.42	0.43
19:AS:22:LEU:HD21	19:AS:28:LYS:N	2.33	0.43
20:AT:43:LEU:O	20:AT:47:GLY:N	2.52	0.43
20:AT:53:LEU:O	20:AT:56:MET:N	2.52	0.43
20:AT:84:LEU:C	20:AT:84:LEU:HD13	2.39	0.43
23:AW:13:C:H42	23:AW:23:G:H1	1.66	0.43
25:AY:90:LEU:HD21	25:AY:104:PRO:HD2	2.01	0.43
26:B0:40:GLN:O	26:B0:57:PHE:HB2	2.18	0.43
27:B1:18:ILE:HA	27:B1:44:PRO:HD2	2.00	0.43
27:B1:72:GLU:O	27:B1:76:ARG:NH2	2.51	0.43
30:B4:17:GLY:O	30:B4:19:GLY:N	2.39	0.43
31:B5:30:LEU:HD23	31:B5:41:PRO:HA	2.00	0.43
34:B8:35:GLN:NE2	34:B8:36:LYS:NZ	2.67	0.43
35:BA:1145:C:H2'	35:BA:1146:C:C6	2.53	0.43
35:BA:1239:G:H2'	35:BA:1240:U:O4'	2.18	0.43
35:BA:1268:A:H2	53:BW:88:ARG:NH2	2.17	0.43
35:BA:1644:C:HO2'	35:BA:1645:G:H5'	1.82	0.43
35:BA:1861:G:H2'	35:BA:1862:G:C8	2.53	0.43
27:B1:41:ARG:NH1	35:BA:189:G:O5'	2.51	0.43
35:BA:1938:A:H4'	35:BA:1939:U:OP2	2.18	0.43
35:BA:1997:G:C2	35:BA:1998:G:N7	2.87	0.43
35:BA:2127:G:O3'	35:BA:2128:C:C4'	2.66	0.43
35:BA:2171:A:O2'	35:BA:2172:U:H6	2.00	0.43
35:BA:2228:G:C6	35:BA:2229:C:C4	3.07	0.43
35:BA:2241:A:C2	35:BA:2242:G:C5	3.06	0.43
35:BA:2314:C:H2'	35:BA:2315:G:C8	2.50	0.43
35:BA:2453:A:O2'	35:BA:2454:G:H5'	2.18	0.43
35:BA:2472:G:H5'	35:BA:2473:U:C5'	2.47	0.43
35:BA:2553:G:H3'	35:BA:2554:U:H5''	2.01	0.43
35:BA:259:G:HO2'	35:BA:621:A:HO2'	1.61	0.43
35:BA:2660:A:H2'	35:BA:2661:G:O4'	2.19	0.43
35:BA:537:C:C2	35:BA:538:G:C8	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:542:C:C4	35:BA:543:C:N4	2.84	0.43
35:BA:685:A:C4	35:BA:689:A:N6	2.86	0.43
35:BA:958:U:OP2	47:BQ:14:ARG:NH1	2.51	0.43
36:BB:63:G:N3	36:BB:63:G:H2'	2.32	0.43
39:BE:65:GLY:O	39:BE:66:HIS:C	2.57	0.43
40:BF:164:ARG:O	40:BF:166:ALA:N	2.51	0.43
40:BF:178:PRO:C	40:BF:180:GLY:N	2.72	0.43
40:BF:53:THR:N	40:BF:56:GLU:HB2	2.17	0.43
41:BG:169:ALA:C	41:BG:171:ALA:N	2.70	0.43
42:BH:12:PRO:O	42:BH:13:LYS:CB	2.67	0.43
43:BI:111:PRO:HA	43:BI:114:LEU:CD1	2.49	0.43
44:BN:120:LEU:CD1	44:BN:120:LEU:C	2.87	0.43
45:BO:105:GLU:O	45:BO:108:GLU:HB2	2.19	0.43
45:BO:23:ARG:HG3	45:BO:24:VAL:N	2.33	0.43
46:BP:111:ARG:HA	46:BP:128:HIS:CE1	2.53	0.43
46:BP:115:LEU:HB3	46:BP:131:SER:OG	2.19	0.43
46:BP:58:THR:O	46:BP:58:THR:HG22	2.19	0.43
47:BQ:16:ARG:HH11	47:BQ:16:ARG:HB3	1.78	0.43
48:BR:44:LEU:CD1	48:BR:48:VAL:HG23	2.49	0.43
48:BR:55:ALA:HA	48:BR:80:PHE:CZ	2.54	0.43
50:BT:107:ASP:CG	50:BT:109:GLU:H	2.21	0.43
51:BU:92:ARG:HG3	51:BU:94:ASN:H	1.83	0.43
52:BV:58:VAL:HG12	52:BV:101:GLY:C	2.38	0.43
54:BX:72:LYS:O	54:BX:73:ARG:CB	2.67	0.43
55:BY:28:LYS:O	55:BY:29:GLU:C	2.57	0.43
55:BY:65:ALA:O	55:BY:67:LEU:N	2.51	0.43
56:BZ:91:LEU:HD23	56:BZ:130:PRO:HG3	2.00	0.43
1:CA:1071:C:H5''	5:CE:49:PRO:HG2	1.99	0.43
1:CA:1163:C:H2'	1:CA:1164:G:C8	2.49	0.43
1:CA:1191:A:H5''	3:CC:4:LYS:HZ2	1.82	0.43
1:CA:1466:C:O2'	1:CA:1467:G:H5'	2.18	0.43
1:CA:318:G:C2	1:CA:319:G:C5	3.06	0.43
1:CA:356:A:C4	1:CA:357:G:C8	3.06	0.43
1:CA:597:G:C6	1:CA:644:G:C6	3.07	0.43
1:CA:662:G:O2'	1:CA:836:G:C5'	2.67	0.43
1:CA:797:C:O2'	1:CA:798:G:H5'	2.19	0.43
1:CA:971:G:H5''	1:CA:972:C:H5''	2.01	0.43
4:CD:138:TYR:CD2	4:CD:138:TYR:C	2.92	0.43
4:CD:97:LEU:CD2	4:CD:97:LEU:O	2.66	0.43
5:CE:64:ARG:HH11	5:CE:64:ARG:HG3	1.83	0.43
6:CF:82:ARG:HA	6:CF:82:ARG:HH11	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:133:GLY:HA2	7:CG:136:LYS:HG2	2.01	0.43
7:CG:42:ILE:HG23	7:CG:117:ALA:HA	2.00	0.43
1:CA:972:C:C4'	10:CJ:57:LYS:CG	2.91	0.43
11:CK:126:ARG:O	11:CK:127:LYS:C	2.56	0.43
11:CK:49:GLY:C	11:CK:50:TYR:HD2	2.22	0.43
12:CL:41:ARG:HG2	12:CL:42:THR:H	1.83	0.43
12:CL:78:GLN:O	12:CL:80:HIS:N	2.51	0.43
16:CP:38:TYR:CE1	16:CP:50:LYS:HB3	2.54	0.43
16:CP:48:TRP:CE3	16:CP:49:LEU:N	2.86	0.43
20:CT:36:LEU:O	20:CT:37:SER:C	2.56	0.43
20:CT:78:ALA:O	20:CT:81:LYS:N	2.51	0.43
1:CA:1328:C:P	21:CU:21:TYR:OH	2.76	0.43
25:CY:128:ALA:O	25:CY:132:ILE:CG1	2.65	0.43
25:CY:137:LEU:O	25:CY:138:ASP:C	2.57	0.43
25:CY:167:GLU:O	25:CY:168:PHE:C	2.56	0.43
25:CY:58:VAL:HG12	25:CY:66:LEU:HD11	1.99	0.43
26:D0:56:ASP:OD2	26:D0:58:THR:OG1	2.36	0.43
31:D5:18:ALA:O	31:D5:19:ARG:C	2.55	0.43
31:D5:9:LYS:NZ	31:D5:9:LYS:HB3	2.34	0.43
35:DA:1042:G:N3	35:DA:1114:G:N2	2.66	0.43
35:DA:1257:C:H2'	35:DA:1258:C:H6	1.84	0.43
35:DA:1495:A:H2	35:DA:1496:A:C4	2.36	0.43
35:DA:171:G:C5	35:DA:172:C:C5	3.07	0.43
35:DA:2189:U:H2'	35:DA:2190:G:C4'	2.49	0.43
35:DA:2225:A:H4'	35:DA:2226:C:C5'	2.47	0.43
35:DA:2246:G:H2'	35:DA:2247:A:H8	1.81	0.43
35:DA:2290:G:H1	35:DA:2342:C:N4	2.17	0.43
35:DA:2330:G:H1	35:DA:2385:C:H42	1.66	0.43
35:DA:2540:C:N4	35:DA:2541:A:C6	2.86	0.43
35:DA:2714:G:C6	35:DA:2715:C:N3	2.86	0.43
35:DA:2805:G:H2'	35:DA:2807:G:C8	2.54	0.43
35:DA:2819:G:H2'	35:DA:2821:A:N7	2.33	0.43
35:DA:2884:U:H2'	35:DA:2885:C:C5'	2.46	0.43
35:DA:532:A:H4'	35:DA:533:G:C1'	2.48	0.43
35:DA:679:C:O2	35:DA:679:C:H2'	2.18	0.43
35:DA:833:U:H2'	35:DA:834:C:H6	1.81	0.43
35:DA:869:G:H2'	35:DA:870:A:O4'	2.19	0.43
35:DA:910:A:N1	35:DA:2277:G:H1'	2.33	0.43
35:DA:98:G:N3	35:DA:98:G:H2'	2.32	0.43
37:DC:65:PRO:HG2	37:DC:189:ILE:CA	2.48	0.43
38:DD:14:ARG:CG	38:DD:14:ARG:NH1	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:8:PRO:HB3	38:DD:14:ARG:HA	1.99	0.43
35:DA:1800:C:OP1	38:DD:264:LYS:NZ	2.51	0.43
38:DD:61:LEU:O	38:DD:63:ARG:NH1	2.51	0.43
41:DG:135:LEU:HB3	41:DG:155:MET:HG2	2.01	0.43
41:DG:18:GLU:HA	41:DG:18:GLU:OE1	2.18	0.43
41:DG:26:GLN:N	41:DG:30:GLU:OE2	2.51	0.43
41:DG:39:ILE:O	41:DG:39:ILE:HG12	2.18	0.43
43:DI:1:MET:HB2	43:DI:21:VAL:O	2.18	0.43
44:DN:120:LEU:CD1	44:DN:120:LEU:C	2.86	0.43
46:DP:95:VAL:HG23	46:DP:125:VAL:CG2	2.48	0.43
49:DS:25:ARG:HD3	49:DS:42:ASP:OD2	2.19	0.43
49:DS:83:LYS:O	49:DS:85:VAL:HG13	2.19	0.43
49:DS:92:TYR:CG	49:DS:93:LYS:N	2.83	0.43
50:DT:115:ARG:HB3	50:DT:116:ALA:H	1.67	0.43
51:DU:74:LEU:HD22	51:DU:78:THR:CG2	2.48	0.43
53:DW:61:ASN:HA	53:DW:61:ASN:HD22	1.55	0.43
53:DW:9:TYR:H	53:DW:9:TYR:HD2	1.66	0.43
54:DX:18:TYR:HA	54:DX:21:PHE:CG	2.53	0.43
55:DY:81:LYS:HA	55:DY:82:PRO:HD3	1.86	0.43
55:DY:84:ARG:O	55:DY:85:VAL:HG22	2.18	0.43
1:AA:1113:C:H6	1:AA:1113:C:O5'	2.02	0.43
1:AA:1157:A:H4'	1:AA:1158:C:O5'	2.18	0.43
1:AA:1230:C:C2'	1:AA:1231:G:H5'	2.49	0.43
1:AA:1459:C:C4	1:AA:1460:A:N7	2.87	0.43
1:AA:1500:A:OP1	1:AA:1505:G:OP1	2.37	0.43
1:AA:429:U:H4'	1:AA:430:A:OP1	2.18	0.43
1:AA:450:G:OP1	1:AA:452:A:OP2	2.36	0.43
1:AA:502:G:C6	1:AA:503:C:C4	3.07	0.43
1:AA:60:A:H4'	1:AA:61:G:O5'	2.19	0.43
1:AA:64:G:H4'	1:AA:66:G:OP1	2.18	0.43
3:AC:104:GLN:NE2	3:AC:105:GLU:H	2.16	0.43
1:AA:1056:U:C5'	3:AC:163:ALA:HB2	2.48	0.43
4:AD:14:ARG:C	4:AD:16:GLY:N	2.72	0.43
9:AI:3:GLN:HA	9:AI:19:LEU:O	2.18	0.43
12:AL:5:PRO:CG	12:AL:10:LEU:HD21	2.49	0.43
12:AL:114:LYS:HE2	12:AL:114:LYS:HB3	1.93	0.43
1:AA:563:A:H2	12:AL:15:ARG:NH1	2.17	0.43
12:AL:6:THR:HG23	12:AL:9:GLN:N	2.12	0.43
13:AM:14:ARG:HB2	13:AM:16:ASP:OD2	2.18	0.43
13:AM:92:HIS:ND1	13:AM:98:VAL:HG21	2.34	0.43
15:AO:85:LEU:HB2	15:AO:87:ILE:HG13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AP:14:ASN:H	16:AP:15:PRO:CD	2.28	0.43
16:AP:53:VAL:HG23	16:AP:54:GLU:N	2.31	0.43
17:AQ:15:MET:HG2	17:AQ:16:GLN:N	2.33	0.43
19:AS:22:LEU:HD13	19:AS:22:LEU:C	2.37	0.43
23:AW:22:A:C6	23:AW:49:C:C2	3.07	0.43
27:B1:23:LYS:N	27:B1:38:SER:O	2.51	0.43
27:B1:89:GLU:O	27:B1:93:GLU:CD	2.56	0.43
32:B6:15:GLU:O	32:B6:16:CYS:O	2.37	0.43
32:B6:28:ARG:HA	32:B6:32:ASN:CB	2.48	0.43
35:BA:1149:G:H2'	35:BA:1150:C:C6	2.54	0.43
35:BA:1778:U:H5	35:BA:1784:A:N3	2.17	0.43
35:BA:1902:C:H5'	38:BD:246:PRO:CD	2.49	0.43
35:BA:198:C:H5'	35:BA:2244:U:OP1	2.19	0.43
35:BA:2415:G:C6	35:BA:2416:C:C4	3.07	0.43
35:BA:2633:G:N2	35:BA:2634:G:H1'	2.33	0.43
35:BA:2670:A:C2	35:BA:2671:A:C8	3.06	0.43
35:BA:402:A:C2'	35:BA:403:U:H5'	2.48	0.43
35:BA:78:A:O2'	35:BA:79:G:H5'	2.19	0.43
35:BA:812:C:O2	35:BA:1250:G:N1	2.51	0.43
38:BD:139:GLY:O	38:BD:140:THR:C	2.56	0.43
38:BD:31:LYS:HA	38:BD:31:LYS:HZ2	1.75	0.43
38:BD:43:ARG:HB3	38:BD:54:ARG:CB	2.49	0.43
39:BE:103:ASP:HA	39:BE:168:MET:HA	2.00	0.43
39:BE:4:ILE:CD1	39:BE:28:ALA:HB1	2.48	0.43
39:BE:47:VAL:CG1	39:BE:49:LEU:HD11	2.48	0.43
40:BF:80:ALA:O	40:BF:83:PHE:HB2	2.18	0.43
42:BH:86:GLU:HA	42:BH:132:ARG:CB	2.49	0.43
43:BI:86:THR:O	43:BI:122:GLU:OE2	2.36	0.43
43:BI:54:GLN:HA	43:BI:57:ARG:HB3	2.00	0.43
43:BI:85:GLU:HB3	43:BI:86:THR:H	1.51	0.43
47:BQ:113:GLN:O	47:BQ:116:GLU:HB3	2.19	0.43
47:BQ:58:PHE:HD1	47:BQ:58:PHE:O	2.02	0.43
47:BQ:68:ILE:N	47:BQ:68:ILE:CD1	2.79	0.43
48:BR:12:ARG:HG3	48:BR:12:ARG:NH1	2.33	0.43
49:BS:51:ALA:HA	49:BS:56:LEU:HD13	2.01	0.43
50:BT:112:ARG:HD3	50:BT:112:ARG:C	2.39	0.43
50:BT:105:LEU:O	50:BT:113:LYS:NZ	2.52	0.43
51:BU:63:VAL:O	51:BU:64:ARG:C	2.57	0.43
1:CA:1361:G:H2'	1:CA:1362:C:C6	2.53	0.43
1:CA:1414:U:O2	1:CA:1487:G:N2	2.51	0.43
1:CA:318:G:H2'	1:CA:319:G:H8	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:408:A:C2	1:CA:409:G:C4	3.06	0.43
1:CA:452:A:C4	1:CA:453:A:C8	3.07	0.43
1:CA:55:A:N7	1:CA:56:U:C4	2.87	0.43
1:CA:690:G:N1	1:CA:691:G:C2	2.87	0.43
1:CA:828:A:N7	1:CA:859:A:C8	2.87	0.43
1:CA:91:C:O5'	1:CA:91:C:H6	2.01	0.43
1:CA:971:G:C4'	1:CA:972:C:H5''	2.43	0.43
3:CC:104:GLN:NE2	3:CC:105:GLU:H	2.16	0.43
4:CD:66:ARG:O	4:CD:69:GLY:N	2.51	0.43
5:CE:101:ILE:HD11	5:CE:119:LEU:HD23	2.01	0.43
5:CE:73:ASN:C	5:CE:73:ASN:HD22	2.21	0.43
7:CG:11:GLN:CG	7:CG:12:LEU:N	2.82	0.43
10:CJ:28:ARG:NH2	10:CJ:34:VAL:O	2.52	0.43
13:CM:51:ALA:O	13:CM:55:ARG:HB2	2.18	0.43
1:CA:668:G:O2'	15:CO:46:HIS:CD2	2.72	0.43
15:CO:87:ILE:O	15:CO:88:ARG:HB2	2.17	0.43
17:CQ:14:LYS:HB2	17:CQ:14:LYS:HZ3	1.83	0.43
17:CQ:68:ARG:HH11	17:CQ:68:ARG:HG2	1.84	0.43
25:CY:107:THR:C	25:CY:111:ARG:NH1	2.72	0.43
25:CY:72:ASP:HB3	25:CY:75:ALA:CB	2.49	0.43
27:D1:74:VAL:C	27:D1:76:ARG:H	2.21	0.43
32:D6:32:ASN:HD22	32:D6:33:LYS:HE3	1.83	0.43
35:DA:1224:C:O2'	35:DA:1225:G:H5'	2.19	0.43
35:DA:1638:C:H2'	35:DA:1639:U:O4'	2.17	0.43
35:DA:1656:C:O2'	35:DA:1657:C:H5'	2.19	0.43
35:DA:1923:U:H2'	35:DA:1924:C:C6	2.54	0.43
35:DA:2006:C:H2'	35:DA:2007:C:C6	2.53	0.43
35:DA:2064:C:H1'	35:DA:2450:A:C5	2.54	0.43
35:DA:2274:A:N6	35:DA:2276:G:C4	2.86	0.43
35:DA:2310:A:N7	41:DG:75:LYS:NZ	2.61	0.43
35:DA:2382:G:C3'	35:DA:2383:G:H5'	2.49	0.43
35:DA:2447:G:H2'	35:DA:2500:U:OP2	2.18	0.43
35:DA:2476:A:N3	35:DA:2476:A:H2'	2.33	0.43
35:DA:2602:A:H4'	35:DA:2603:G:O5'	2.19	0.43
35:DA:304:G:C2	35:DA:305:U:C2	3.07	0.43
35:DA:475:U:O4'	35:DA:509:C:C5	2.72	0.43
35:DA:677:A:N6	35:DA:802:A:N7	2.67	0.43
35:DA:730:C:O2'	35:DA:731:C:H5'	2.18	0.43
35:DA:792:G:N7	35:DA:2440:C:H1'	2.34	0.43
38:DD:213:ARG:C	38:DD:215:LEU:N	2.72	0.43
38:DD:20:ASP:OD1	38:DD:21:PHE:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DD:240:ALA:HA	38:DD:241:PRO:HD3	1.85	0.43
38:DD:27:THR:HG21	38:DD:83:GLU:CG	2.30	0.43
39:DE:188:VAL:HG13	39:DE:188:VAL:O	2.19	0.43
39:DE:63:LEU:O	39:DE:65:GLY:N	2.52	0.43
40:DF:160:ASN:HD21	40:DF:162:LEU:HB2	1.83	0.43
41:DG:5:VAL:CG1	41:DG:6:ALA:N	2.61	0.43
42:DH:160:LYS:HB3	42:DH:161:GLY:H	1.72	0.43
43:DI:60:GLU:C	43:DI:62:LYS:H	2.22	0.43
43:DI:81:VAL:O	43:DI:83:ALA:N	2.52	0.43
44:DN:3:THR:C	44:DN:5:VAL:N	2.71	0.43
46:DP:62:LEU:H	46:DP:62:LEU:HD12	1.80	0.43
47:DQ:20:ALA:HA	47:DQ:98:LYS:CB	2.45	0.43
49:DS:106:ARG:HD2	49:DS:106:ARG:O	2.19	0.43
49:DS:87:PHE:CE2	49:DS:88:ASP:O	2.72	0.43
54:DX:59:VAL:HG22	54:DX:74:PRO:O	2.19	0.43
54:DX:65:ARG:NE	54:DX:66:LEU:N	2.65	0.43
54:DX:76:ARG:O	54:DX:77:LYS:CB	2.65	0.43
55:DY:84:ARG:C	55:DY:85:VAL:CG2	2.86	0.43
56:DZ:97:GLU:HB3	56:DZ:125:LEU:CD2	2.37	0.43
1:AA:1131:G:C2	1:AA:1132:C:N4	2.87	0.43
1:AA:1361:G:H2'	1:AA:1362:C:C6	2.54	0.43
1:AA:922:G:N3	1:AA:1396:A:C2	2.87	0.43
1:AA:1434:A:H2'	1:AA:1435:G:H8	1.83	0.43
1:AA:1442(A):G:H5'	1:AA:1442(B):A:C2	2.53	0.43
1:AA:1496:C:H2'	1:AA:1497:G:O4'	2.19	0.43
1:AA:1506:U:OP1	24:AX:15:G:OP1	2.37	0.43
1:AA:15:G:O3'	5:AE:24:ARG:NH2	2.52	0.43
1:AA:356:A:C4	1:AA:357:G:C8	3.07	0.43
1:AA:35:G:C4	1:AA:36:C:C5	3.06	0.43
1:AA:515:G:N2	1:AA:537:G:C4	2.87	0.43
1:AA:61:G:H2'	1:AA:62:U:O4'	2.18	0.43
1:AA:706:A:N7	1:AA:707:C:H5	2.17	0.43
1:AA:680:C:N3	1:AA:711:G:C2	2.86	0.43
1:AA:725:G:O2'	1:AA:726:C:H5'	2.19	0.43
1:AA:825:G:O2'	1:AA:826:C:H5'	2.18	0.43
1:AA:920:U:C2	1:AA:921:U:C5	3.06	0.43
1:AA:932:C:H5''	7:AG:3:ARG:CD	2.37	0.43
2:AB:105:PHE:HA	2:AB:108:ILE:CG2	2.48	0.43
2:AB:159:PRO:HB2	2:AB:162:ILE:HG22	2.01	0.43
2:AB:166:ASP:O	2:AB:167:PRO:O	2.37	0.43
2:AB:84:GLU:HB3	2:AB:219:VAL:CG2	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:11:ARG:HG2	3:AC:11:ARG:NH1	2.32	0.43
4:AD:172:PRO:C	4:AD:187:ARG:HH12	2.22	0.43
4:AD:45:GLN:C	4:AD:46:LYS:HG3	2.39	0.43
4:AD:98:GLU:C	4:AD:100:ARG:N	2.72	0.43
6:AF:70:ASP:O	6:AF:71:ARG:C	2.57	0.43
7:AG:16:LEU:HD11	9:AI:42:ARG:HA	2.00	0.43
8:AH:16:ALA:HB1	8:AH:63:LEU:HD21	2.01	0.43
9:AI:28:VAL:O	9:AI:30:GLY:N	2.51	0.43
11:AK:99:GLN:O	11:AK:101:SER:N	2.52	0.43
13:AM:40:ASN:HA	13:AM:41:PRO:HD3	1.79	0.43
13:AM:9:ILE:N	13:AM:9:ILE:CD1	2.82	0.43
16:AP:27:LYS:H	16:AP:27:LYS:CD	2.28	0.43
18:AR:74:ARG:CZ	18:AR:81:PHE:HA	2.49	0.43
19:AS:5:LEU:H	19:AS:6:LYS:NZ	2.17	0.43
21:AU:19:GLY:O	21:AU:22:ARG:O	2.36	0.43
25:AY:154:THR:O	25:AY:157:ALA:HB3	2.19	0.43
26:B0:25:ARG:CB	26:B0:37:LEU:HD23	2.46	0.43
27:B1:34:THR:O	27:B1:35:THR:HG23	2.18	0.43
27:B1:62:VAL:O	27:B1:63:ALA:O	2.37	0.43
28:B2:29:LYS:O	28:B2:32:LEU:N	2.52	0.43
28:B2:47:ASN:CG	28:B2:48:HIS:N	2.71	0.43
34:B8:46:ARG:NH1	34:B8:46:ARG:CB	2.60	0.43
35:BA:1051:G:C2	35:BA:1052:C:N4	2.87	0.43
35:BA:1195:G:C2	35:BA:1196:C:C4	3.06	0.43
35:BA:1210:A:C8	35:BA:1237:A:N6	2.87	0.43
35:BA:127:A:H5''	35:BA:128:C:C6	2.54	0.43
35:BA:1298:C:C2'	35:BA:1299:G:C8	3.02	0.43
35:BA:1301:A:C2	35:BA:1303:G:C6	3.07	0.43
35:BA:1358:G:N7	35:BA:1371:G:C6	2.86	0.43
35:BA:1445(A):C:H2'	35:BA:1446:C:C6	2.52	0.43
35:BA:1444:G:O6	35:BA:1466:G:C6	2.72	0.43
35:BA:1466:G:N2	35:BA:1547:C:C2	2.87	0.43
35:BA:1853:A:H2'	35:BA:1854:A:O4'	2.19	0.43
35:BA:2102:U:O4'	35:BA:2102:U:O2	2.36	0.43
35:BA:2447:G:H2'	35:BA:2500:U:OP2	2.18	0.43
35:BA:2591:C:H6	35:BA:2591:C:O5'	2.02	0.43
35:BA:259:G:C2	35:BA:260:G:C8	3.07	0.43
35:BA:366:C:H5''	35:BA:403:U:N3	2.33	0.43
35:BA:392:C:H5''	35:BA:409:C:H5''	2.01	0.43
35:BA:579:G:C6	35:BA:580:C:N4	2.87	0.43
35:BA:724:U:C2'	35:BA:725:G:H5'	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:835:A:O2'	35:BA:836:G:H5'	2.19	0.43
36:BB:7:G:H4'	49:BS:29:PHE:HE2	1.80	0.43
35:BA:1788:C:OP1	38:BD:222:ARG:NH2	2.52	0.43
38:BD:231:HIS:NE2	38:BD:249:PRO:HG3	2.33	0.43
41:BG:110:ALA:O	41:BG:113:ARG:N	2.52	0.43
41:BG:173:LEU:HA	41:BG:176:LEU:CD1	2.27	0.43
44:BN:62:VAL:HG21	44:BN:66:LYS:HB2	2.00	0.43
35:BA:1245:G:OP1	46:BP:16:ARG:NE	2.52	0.43
48:BR:6:SER:HA	48:BR:8:ARG:HH21	1.79	0.43
49:BS:31:SER:CB	49:BS:34:HIS:O	2.67	0.43
49:BS:51:ALA:HA	49:BS:56:LEU:CD1	2.48	0.43
45:BO:78:ARG:HB3	50:BT:73:GLU:CG	2.48	0.43
50:BT:82:LEU:O	50:BT:83:ILE:HG12	2.19	0.43
56:BZ:47:VAL:O	56:BZ:51:ALA:CB	2.66	0.43
1:CA:1011:G:H2'	1:CA:1012:U:H5'	2.01	0.43
1:CA:1132:C:C2'	1:CA:1133:G:H5'	2.49	0.43
1:CA:1131:G:C2	1:CA:1132:C:N4	2.87	0.43
1:CA:1340:A:P	23:CW:36:A:H4'	2.58	0.43
1:CA:1479:C:H2'	1:CA:1480:G:H8	1.83	0.43
1:CA:1502:A:H2	1:CA:1505:G:H22	1.64	0.43
1:CA:21:G:H2'	1:CA:22:G:C8	2.54	0.43
1:CA:245:C:O2	1:CA:283:C:N3	2.51	0.43
1:CA:357:G:O2'	1:CA:358:U:H5'	2.19	0.43
1:CA:630:G:H2'	1:CA:631:G:H5'	2.00	0.43
1:CA:781:A:H4'	1:CA:1522:U:O2'	2.19	0.43
1:CA:786:G:C2	1:CA:797:C:O2	2.72	0.43
1:CA:821:G:O2'	1:CA:822:C:H5'	2.18	0.43
1:CA:984:C:H2'	1:CA:985:C:C6	2.54	0.43
2:CB:164:VAL:O	2:CB:186:ALA:CB	2.67	0.43
3:CC:190:ARG:O	3:CC:191:THR:O	2.36	0.43
4:CD:59:ARG:HA	4:CD:59:ARG:CZ	2.49	0.43
5:CE:48:ALA:O	5:CE:50:GLU:N	2.52	0.43
7:CG:26:PHE:O	7:CG:30:ILE:HG12	2.19	0.43
7:CG:57:GLU:C	7:CG:59:LEU:H	2.22	0.43
7:CG:65:ALA:HB2	7:CG:124:LEU:O	2.19	0.43
8:CH:114:THR:CG2	8:CH:117:GLY:O	2.67	0.43
8:CH:5:PRO:HA	8:CH:8:ASP:HB3	2.00	0.43
1:CA:1368:G:OP2	9:CI:112:LYS:HE3	2.18	0.43
1:CA:973:G:P	10:CJ:57:LYS:NZ	2.92	0.43
1:CA:684:A:C2	11:CK:39:PRO:HG2	2.54	0.43
11:CK:99:GLN:O	11:CK:101:SER:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:126:LYS:HA	12:CL:126:LYS:HD2	1.76	0.43
1:CA:563:A:H2	12:CL:15:ARG:NH1	2.16	0.43
13:CM:33:ALA:HA	13:CM:59:TYR:CE2	2.54	0.43
14:CN:3:ARG:O	14:CN:7:ILE:HG23	2.18	0.43
14:CN:6:LEU:C	14:CN:8:GLU:N	2.72	0.43
16:CP:18:ARG:HD3	16:CP:35:LYS:CD	2.46	0.43
16:CP:55:ARG:O	16:CP:58:TYR:HB3	2.19	0.43
17:CQ:10:VAL:CG1	17:CQ:53:LEU:HA	2.49	0.43
19:CS:10:PHE:CE2	19:CS:70:LYS:NZ	2.87	0.43
19:CS:36:ARG:HH22	19:CS:75:ALA:CB	2.17	0.43
20:CT:82:SER:O	20:CT:86:ARG:HB2	2.19	0.43
25:CY:68:VAL:C	25:CY:98:ALA:HA	2.38	0.43
27:D1:18:ILE:HG12	27:D1:43:TYR:HD1	1.84	0.43
29:D3:22:ALA:O	29:D3:26:LEU:HG	2.19	0.43
34:D8:2:PRO:HA	35:DA:591:C:O2	2.18	0.43
35:DA:1208:C:C4	35:DA:1209:G:N7	2.87	0.43
35:DA:1210:A:C8	35:DA:1237:A:N6	2.87	0.43
35:DA:1374:G:H2'	35:DA:1375:C:H6	1.83	0.43
35:DA:154(A):C:O2	35:DA:154(A):C:O4'	2.33	0.43
35:DA:1560:G:OP1	35:DA:1561:G:OP2	2.37	0.43
35:DA:1569:A:O2'	35:DA:1570:A:H5'	2.19	0.43
35:DA:1605:C:H2'	35:DA:1606:G:O4'	2.18	0.43
35:DA:1683:C:H2'	35:DA:1684:C:C6	2.53	0.43
35:DA:1826:G:H2'	35:DA:1827:C:H6	1.84	0.43
35:DA:2206:G:N3	35:DA:2207:G:H5'	2.33	0.43
35:DA:2543:G:O2'	35:DA:2544:G:H5'	2.19	0.43
35:DA:2713:A:H3'	35:DA:2714:G:H5'	2.00	0.43
35:DA:2864:G:C8	35:DA:2864:G:H5'	2.47	0.43
35:DA:483:A:N7	35:DA:497:A:H2	2.16	0.43
35:DA:491:G:N2	35:DA:492:A:H1'	2.34	0.43
37:DC:184:LYS:C	37:DC:186:ALA:N	2.72	0.43
38:DD:132:PRO:HB3	38:DD:188:GLU:O	2.19	0.43
38:DD:3:VAL:HG12	38:DD:4:LYS:N	2.33	0.43
41:DG:103:LEU:O	41:DG:107:LEU:HB2	2.18	0.43
41:DG:116:ASP:CG	41:DG:117:PHE:H	2.22	0.43
42:DH:136:ILE:CD1	42:DH:136:ILE:N	2.80	0.43
42:DH:83:TYR:HD1	42:DH:84:SER:N	2.16	0.43
43:DI:132:PRO:O	43:DI:133:HIS:O	2.36	0.43
44:DN:128:HIS:HD2	44:DN:131:GLN:CB	2.31	0.43
44:DN:38:HIS:O	44:DN:40:PRO:HD3	2.18	0.43
44:DN:46:VAL:HG11	44:DN:48:MET:CG	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:66:LYS:HE3	44:DN:66:LYS:HA	1.99	0.43
45:DO:10:VAL:HG21	45:DO:16:ALA:O	2.18	0.43
46:DP:65:ARG:O	46:DP:66:GLY:C	2.57	0.43
48:DR:2:ARG:NE	48:DR:5:LYS:HZ2	2.14	0.43
49:DS:28:VAL:H	49:DS:89:ARG:CB	2.30	0.43
49:DS:74:ALA:O	49:DS:75:GLU:C	2.57	0.43
55:DY:37:VAL:CG2	55:DY:38:ILE:N	2.71	0.43
55:DY:43:ASN:CG	55:DY:64:GLU:HA	2.38	0.43
56:DZ:147:GLY:O	56:DZ:148:ASP:CB	2.66	0.43
56:DZ:153:SER:C	56:DZ:154:ASP:OD2	2.57	0.43
56:DZ:69:THR:CG2	56:DZ:90:VAL:HG22	2.29	0.43
1:AA:1050:G:N2	1:AA:1209:C:H1'	2.34	0.43
1:AA:1157:A:H1'	1:AA:1181:G:N2	2.34	0.43
1:AA:1237:C:C4'	1:AA:1334:G:N2	2.82	0.43
1:AA:129(A):G:N2	1:AA:189(F):U:H5''	2.34	0.43
1:AA:346:G:N3	1:AA:346:G:H2'	2.33	0.43
1:AA:106:C:O2'	1:AA:379:C:OP1	2.37	0.43
1:AA:452:A:H1'	16:AP:72:ARG:NH1	2.34	0.43
1:AA:783:C:N4	1:AA:784:C:H41	2.16	0.43
1:AA:799:G:H2'	1:AA:800:G:H5'	2.01	0.43
1:AA:919:A:H2'	1:AA:920:U:O5'	2.19	0.43
3:AC:180:ALA:O	3:AC:181:ASN:O	2.37	0.43
3:AC:73:PRO:O	3:AC:76:VAL:N	2.52	0.43
4:AD:126:ILE:CG2	4:AD:127:THR:H	2.30	0.43
4:AD:17:VAL:CG1	4:AD:18:LYS:N	2.80	0.43
5:AE:64:ARG:HG3	5:AE:64:ARG:HH11	1.84	0.43
8:AH:86:ILE:HG21	8:AH:133:LEU:CD2	2.49	0.43
9:AI:28:VAL:HG22	9:AI:63:ILE:HB	1.99	0.43
9:AI:24:GLY:CA	9:AI:60:ASP:HA	2.49	0.43
9:AI:75:ASP:O	9:AI:78:LYS:NZ	2.44	0.43
11:AK:51:LYS:HA	11:AK:55:LYS:HG3	2.01	0.43
11:AK:88:GLY:C	11:AK:91:ARG:HB2	2.38	0.43
14:AN:47:LEU:O	14:AN:51:GLY:N	2.52	0.43
15:AO:9:GLN:HB3	15:AO:13:GLN:NE2	2.17	0.43
16:AP:34:GLU:OE1	16:AP:36:ILE:HG23	2.19	0.43
20:AT:24:LEU:O	20:AT:25:ARG:C	2.57	0.43
23:AW:16:C:H6	23:AW:60:A:H2	1.67	0.43
25:AY:96:GLY:C	25:AY:98:ALA:H	2.23	0.43
26:B0:40:GLN:HG3	26:B0:57:PHE:HB3	2.00	0.43
27:B1:48:LYS:O	27:B1:49:VAL:CG2	2.65	0.43
31:B5:54:GLY:O	31:B5:56:LYS:NZ	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:34:ARG:HE	33:B7:39:ARG:HD2	1.83	0.43
33:B7:41:ARG:O	33:B7:42:LEU:C	2.54	0.43
35:BA:1231:G:H2'	35:BA:1232:G:C8	2.54	0.43
35:BA:1260:G:H2'	35:BA:1261:C:C6	2.54	0.43
35:BA:1354:A:H2'	35:BA:1355:G:H5'	2.00	0.43
35:BA:1400:G:H2'	35:BA:1401:G:H8	1.84	0.43
35:BA:1403:C:C2'	35:BA:1404:C:O5'	2.67	0.43
35:BA:1613:G:C2'	35:BA:1617:C:N4	2.81	0.43
35:BA:1658:C:H2'	35:BA:1659:U:H6	1.79	0.43
35:BA:1664:A:H61	35:BA:1996:C:H42	1.67	0.43
35:BA:1680:U:O2	35:BA:1763:G:H3'	2.18	0.43
35:BA:2320:A:H2'	35:BA:2320:A:N3	2.34	0.43
35:BA:2369:A:H2'	35:BA:2370:G:H8	1.84	0.43
35:BA:2447:G:C6	35:BA:2501:C:C2	3.07	0.43
35:BA:2550:G:H2'	35:BA:2551:C:C6	2.54	0.43
35:BA:2678:C:H2'	35:BA:2678:C:O2	2.17	0.43
35:BA:2679:A:H2'	35:BA:2680:C:C6	2.54	0.43
35:BA:2711:A:OP2	35:BA:2712(A):A:OP2	2.37	0.43
35:BA:2733:A:H2'	35:BA:2734:A:O4'	2.18	0.43
35:BA:418:G:O2'	35:BA:419:C:H5'	2.19	0.43
35:BA:534:U:O3'	51:BU:46:ALA:CB	2.67	0.43
35:BA:556:G:C5	35:BA:557:U:C4	3.07	0.43
35:BA:614(B):G:C4	40:BF:44:ARG:NH2	2.86	0.43
35:BA:863:A:OP1	47:BQ:21:THR:HB	2.18	0.43
35:BA:923:C:H2'	35:BA:924:C:C6	2.53	0.43
35:BA:936:C:H2'	35:BA:937:U:C6	2.54	0.43
36:BB:64:C:H2'	36:BB:65:C:C6	2.54	0.43
38:BD:95:LEU:HD11	38:BD:105:ILE:HG22	2.00	0.43
35:BA:2052:G:P	39:BE:141:ILE:HD11	2.59	0.43
39:BE:23:VAL:CA	39:BE:184:VAL:O	2.66	0.43
42:BH:127:GLU:CB	42:BH:130:ARG:HB3	2.49	0.43
44:BN:56:ASN:CA	44:BN:125:GLY:H	2.29	0.43
44:BN:7:LYS:O	44:BN:8:GLN:C	2.58	0.43
45:BO:87:ILE:HG23	45:BO:88:ASN:O	2.19	0.43
46:BP:29:LYS:HD2	46:BP:29:LYS:N	2.33	0.43
46:BP:57:THR:O	46:BP:57:THR:HG22	2.19	0.43
35:BA:389:G:N1	46:BP:71:VAL:CG1	2.80	0.43
47:BQ:57:HIS:ND1	47:BQ:58:PHE:N	2.66	0.43
48:BR:10:LEU:HD22	48:BR:17:ARG:CD	2.31	0.43
49:BS:106:ARG:HD2	49:BS:107:GLU:O	2.19	0.43
49:BS:58:LEU:CD2	49:BS:65:VAL:HG13	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BO:71:ARG:HH12	50:BT:74:ARG:HH22	1.66	0.43
50:BT:84:GLN:O	50:BT:84:GLN:HG3	2.19	0.43
51:BU:49:HIS:O	51:BU:52:ARG:N	2.52	0.43
51:BU:51:LYS:CA	51:BU:51:LYS:HE2	2.49	0.43
35:BA:1160:G:H22	52:BV:10:LYS:HE3	1.84	0.43
51:BU:43:GLY:HA3	52:BV:76:LYS:HB2	2.00	0.43
53:BW:69:LEU:N	53:BW:69:LEU:HD12	2.32	0.43
54:BX:89:ILE:O	54:BX:89:ILE:CG2	2.66	0.43
55:BY:31:LEU:CG	55:BY:34:LYS:HB2	2.49	0.43
56:BZ:11:GLU:HB2	56:BZ:13:GLU:CG	2.47	0.43
56:BZ:27:VAL:N	56:BZ:86:VAL:O	2.48	0.43
56:BZ:6:LYS:HB3	56:BZ:8:TYR:CE1	2.53	0.43
1:CA:1079:G:C2	1:CA:1080:A:C5	3.06	0.43
1:CA:107:G:H2'	1:CA:108:G:H5'	1.99	0.43
1:CA:1157:A:C1'	1:CA:1181:G:N2	2.81	0.43
1:CA:260:G:H2'	1:CA:261:U:C6	2.54	0.43
1:CA:59:A:C2	1:CA:331:G:C2	3.07	0.43
1:CA:338:A:C4	1:CA:339:C:C5	3.07	0.43
1:CA:32:A:C2	1:CA:33:A:C4	3.07	0.43
1:CA:680:C:N3	1:CA:711:G:C2	2.87	0.43
1:CA:729:A:H2'	1:CA:730:G:C8	2.53	0.43
2:CB:173:ALA:O	2:CB:174:VAL:C	2.55	0.43
4:CD:138:TYR:C	4:CD:138:TYR:HD2	2.23	0.43
8:CH:26:VAL:CG1	8:CH:59:LEU:HB2	2.48	0.43
9:CI:27:THR:CG2	9:CI:28:VAL:N	2.82	0.43
9:CI:98:PRO:O	9:CI:100:GLY:N	2.52	0.43
11:CK:115:PRO:C	11:CK:116:HIS:ND1	2.72	0.43
11:CK:124:LYS:HD2	11:CK:125:PHE:CE1	2.53	0.43
1:CA:523:A:N6	12:CL:53:ARG:NH2	2.66	0.43
12:CL:6:THR:H	12:CL:9:GLN:NE2	2.17	0.43
13:CM:14:ARG:HB2	13:CM:16:ASP:OD2	2.18	0.43
1:CA:1219:U:P	14:CN:19:ARG:HH22	2.42	0.43
17:CQ:63:ARG:HG2	17:CQ:64:PRO:HD2	2.01	0.43
18:CR:43:PHE:N	18:CR:43:PHE:CD1	2.78	0.43
19:CS:24:ALA:O	19:CS:25:LYS:HB2	2.19	0.43
20:CT:64:ASP:OD1	20:CT:81:LYS:HD2	2.19	0.43
23:CW:33:C:O2	23:CW:37:U:H5	2.01	0.43
24:CX:15:G:H5''	24:CX:16:U:OP1	2.19	0.43
25:CY:52:LEU:O	25:CY:53:ASN:C	2.56	0.43
25:CY:5:GLU:O	25:CY:6:LEU:C	2.56	0.43
25:CY:84:ARG:C	25:CY:86:SER:N	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:D3:22:ALA:HB1	29:D3:46:ASN:HD22	1.84	0.43
35:DA:1142(A):A:C4	35:DA:1144:G:C8	3.07	0.43
35:DA:1153:C:C4	35:DA:1154:G:C2	3.07	0.43
35:DA:1315:C:H42	35:DA:1337:G:H1	1.66	0.43
35:DA:1338:G:N3	35:DA:1338:G:H2'	2.34	0.43
35:DA:1491:G:H5'	38:DD:99:ASP:OD1	2.19	0.43
35:DA:1544:A:H2'	35:DA:1545:A:OP2	2.19	0.43
35:DA:1560:G:C2	35:DA:1561:G:C4	3.07	0.43
35:DA:1635:G:H2'	35:DA:1636:C:H6	1.84	0.43
35:DA:1720:U:H2'	35:DA:1721:G:C5'	2.49	0.43
35:DA:2019:A:C6	35:DA:2020:A:C5	3.06	0.43
35:DA:2319:G:C4	35:DA:2320:A:C6	3.06	0.43
35:DA:2700:C:H2'	35:DA:2701:C:H6	1.83	0.43
35:DA:271(P):C:O2'	35:DA:271(Q):G:H5'	2.19	0.43
35:DA:2733:A:H2'	35:DA:2734:A:O4'	2.18	0.43
35:DA:431:U:C2'	35:DA:432:A:H5'	2.47	0.43
36:DB:110:G:H2'	36:DB:111:G:H8	1.83	0.43
38:DD:66:ASP:OD2	38:DD:103:ARG:NH2	2.51	0.43
38:DD:70:TRP:CZ2	38:DD:150:LYS:HA	2.54	0.43
39:DE:2:LYS:HD3	39:DE:95:ILE:HG22	2.01	0.43
40:DF:102:PRO:O	40:DF:104:LYS:N	2.51	0.43
40:DF:136:THR:O	40:DF:137:LYS:C	2.57	0.43
41:DG:109:VAL:C	41:DG:111:LEU:N	2.68	0.43
41:DG:130:ASN:HB3	41:DG:159:VAL:O	2.18	0.43
42:DH:43:VAL:HB	42:DH:51:ARG:O	2.18	0.43
42:DH:44:VAL:HG12	42:DH:45:VAL:HG23	2.00	0.43
42:DH:79:VAL:O	42:DH:81:GLU:N	2.48	0.43
42:DH:89:ILE:HG12	42:DH:90:LYS:H	1.81	0.43
44:DN:98:VAL:O	44:DN:102:ALA:CB	2.67	0.43
45:DO:105:GLU:O	45:DO:108:GLU:HB2	2.18	0.43
45:DO:88:ASN:HB3	45:DO:92:GLU:O	2.19	0.43
47:DQ:125:LEU:HB3	47:DQ:127:ILE:CD1	2.49	0.43
48:DR:9:LYS:HZ1	48:DR:39:PRO:HA	1.82	0.43
50:DT:48:ILE:HG22	50:DT:49:VAL:N	2.34	0.43
52:DV:83:ARG:O	52:DV:84:LYS:CD	2.67	0.43
52:DV:88:ARG:CG	52:DV:88:ARG:HH11	2.32	0.43
53:DW:5:ALA:CB	53:DW:105:VAL:HB	2.48	0.43
53:DW:4:LYS:HA	53:DW:106:ILE:HA	2.00	0.43
55:DY:31:LEU:HB2	55:DY:36:ALA:N	2.26	0.43
55:DY:31:LEU:HD12	55:DY:34:LYS:H	1.80	0.43
56:DZ:151:HIS:O	56:DZ:171:ILE:CD1	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:151:HIS:O	56:DZ:171:ILE:HD13	2.19	0.43
56:DZ:5:LEU:HD21	56:DZ:43:GLU:HB3	2.01	0.43
1:AA:1105:A:C2	1:AA:1106:G:N7	2.87	0.43
1:AA:1126:U:O2'	1:AA:1127:G:H5'	2.19	0.43
1:AA:1296:C:H3'	1:AA:1297:C:C6	2.51	0.43
1:AA:1441:G:O3'	1:AA:1442:G:C8	2.71	0.43
1:AA:320:C:O4'	1:AA:1434:A:H2	2.02	0.43
1:AA:33:A:H2'	1:AA:34:C:H6	1.83	0.43
1:AA:369:C:OP2	1:AA:388:G:N2	2.46	0.43
1:AA:38:G:C4	1:AA:397:A:N1	2.87	0.43
1:AA:435:C:N3	1:AA:436:C:C5	2.87	0.43
1:AA:585:G:O2'	1:AA:879:C:OP1	2.37	0.43
1:AA:663:A:H2'	1:AA:664:G:C8	2.53	0.43
1:AA:696:A:C4	1:AA:697:U:C5	3.07	0.43
1:AA:832:C:O2'	1:AA:833:U:P	2.76	0.43
1:AA:867:G:N2	1:AA:868:C:C2	2.87	0.43
1:AA:925:G:H1	1:AA:1391:U:H3	1.65	0.43
2:AB:211:ILE:O	2:AB:215:LEU:N	2.47	0.43
2:AB:33:TYR:HB2	2:AB:41:ILE:O	2.19	0.43
3:AC:150:LYS:HB3	3:AC:201:TYR:CB	2.41	0.43
3:AC:173:VAL:HG12	3:AC:175:LEU:HG	2.01	0.43
3:AC:8:ILE:C	3:AC:10:PHE:N	2.72	0.43
4:AD:102:ASP:HB3	4:AD:136:PRO:HB3	2.00	0.43
8:AH:122:ARG:NH1	8:AH:122:ARG:CB	2.82	0.43
8:AH:20:TYR:HA	8:AH:65:TYR:CZ	2.50	0.43
9:AI:126:SER:O	9:AI:128:ARG:HD3	2.19	0.43
10:AJ:56:HIS:O	10:AJ:58:ASP:O	2.36	0.43
10:AJ:5:ARG:HG3	10:AJ:73:ASP:CG	2.39	0.43
10:AJ:8:LEU:HG	10:AJ:96:ILE:CG2	2.49	0.43
12:AL:117:ARG:HD2	12:AL:122:THR:HB	2.00	0.43
14:AN:24:CYS:SG	14:AN:39:LEU:HA	2.59	0.43
10:AJ:50:ILE:HD11	14:AN:41:ARG:HH11	1.84	0.43
18:AR:36:ASN:HD21	18:AR:40:LEU:HD21	1.84	0.43
18:AR:53:ARG:HH22	18:AR:59:SER:C	2.20	0.43
19:AS:43:GLU:O	19:AS:43:GLU:OE1	2.36	0.43
25:AY:32:ARG:NH2	25:AY:88:LEU:HG	2.34	0.43
27:B1:11:ARG:CB	27:B1:12:PRO:CD	2.97	0.43
27:B1:87:PRO:O	27:B1:90:ILE:N	2.34	0.43
27:B1:73:LEU:HD21	27:B1:94:LEU:HB2	1.99	0.43
31:B5:42:PRO:HB2	35:BA:2815:C:HO2'	1.80	0.43
31:B5:51:TYR:HB3	31:B5:52:TYR:H	1.35	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:29:LYS:HZ3	33:B7:32:LYS:HZ2	1.67	0.43
34:B8:39:LYS:HG2	34:B8:42:ARG:HH12	1.80	0.43
35:BA:13:A:H2	35:BA:526:A:C5	2.35	0.43
35:BA:1488:G:N3	35:BA:1488:G:H2'	2.34	0.43
35:BA:1504:C:O2'	35:BA:1505:C:H5'	2.19	0.43
35:BA:1525:G:H2'	35:BA:1526:G:H8	1.83	0.43
35:BA:1578:U:O2	35:BA:1578:U:H2'	2.19	0.43
35:BA:1942:C:C4	35:BA:1943:U:C4	3.07	0.43
35:BA:1992:G:O2'	35:BA:1993:U:P	2.76	0.43
35:BA:1266:G:N2	35:BA:2012:G:C4	2.86	0.43
35:BA:2020:A:C2	35:BA:2035:G:N1	2.87	0.43
35:BA:2083:G:O2'	35:BA:2084:C:H5'	2.19	0.43
35:BA:2205:C:H1'	35:BA:2220:G:N2	2.34	0.43
35:BA:224:G:C6	35:BA:225:A:C5	3.07	0.43
35:BA:2319:G:C4	35:BA:2320:A:C6	3.07	0.43
35:BA:1983:C:H4'	35:BA:2606:C:H4'	2.01	0.43
35:BA:2640:G:C8	35:BA:2640:G:H5'	2.54	0.43
35:BA:270:A:OP2	35:BA:271(X):G:N2	2.47	0.43
35:BA:753:C:H6	35:BA:753:C:O5'	2.02	0.43
35:BA:826:U:H3'	35:BA:828:U:C6	2.54	0.43
35:BA:956:G:H5'	35:BA:957:A:OP2	2.19	0.43
38:BD:19:ALA:O	38:BD:21:PHE:HE1	2.01	0.43
38:BD:57:GLY:O	38:BD:58:HIS:O	2.37	0.43
39:BE:110:GLY:HA2	39:BE:161:GLY:CA	2.30	0.43
39:BE:201:THR:CG2	39:BE:203:LYS:HB3	2.49	0.43
39:BE:9:VAL:CG1	39:BE:25:VAL:HG12	2.49	0.43
39:BE:72:VAL:O	39:BE:73:GLU:O	2.37	0.43
40:BF:192:LEU:C	40:BF:192:LEU:HD23	2.39	0.43
41:BG:18:GLU:O	41:BG:22:ARG:HB2	2.19	0.43
41:BG:36:LYS:HB3	41:BG:36:LYS:HE2	1.90	0.43
41:BG:85:GLY:C	41:BG:87:PRO:HD2	2.39	0.43
44:BN:137:LYS:CG	44:BN:138:LEU:H	2.27	0.43
44:BN:17:ASP:OD2	44:BN:56:ASN:HB3	2.19	0.43
44:BN:87:LEU:O	44:BN:90:MET:HB2	2.19	0.43
45:BO:119:PRO:O	45:BO:120:GLU:HB2	2.17	0.43
46:BP:14:LYS:O	46:BP:15:ARG:CB	2.67	0.43
46:BP:83:VAL:O	46:BP:83:VAL:HG13	2.19	0.43
47:BQ:131:ILE:HD13	47:BQ:131:ILE:H	1.80	0.43
47:BQ:22:LYS:NZ	47:BQ:22:LYS:CA	2.81	0.43
48:BR:53:HIS:C	48:BR:56:LYS:HB2	2.38	0.43
49:BS:19:LYS:HB3	49:BS:20:ARG:NH2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BS:74:ALA:O	49:BS:75:GLU:C	2.58	0.43
49:BS:85:VAL:CG2	49:BS:86:ALA:H	2.17	0.43
50:BT:53:ARG:O	50:BT:53:ARG:HG2	2.19	0.43
50:BT:65:LYS:HE3	50:BT:66:VAL:H	1.83	0.43
51:BU:74:LEU:HD22	51:BU:78:THR:CG2	2.49	0.43
51:BU:83:LEU:HD12	51:BU:113:ALA:HB2	2.00	0.43
56:BZ:24:LEU:HB2	56:BZ:41:LEU:CD2	2.49	0.43
47:BQ:137:TYR:HE2	56:BZ:76:LEU:CD2	2.32	0.43
1:CA:1276:G:C2'	1:CA:1277:C:H5'	2.49	0.43
1:CA:1416:G:C6	1:CA:1485:U:O2	2.72	0.43
1:CA:277:C:P	17:CQ:41:LYS:HZ1	2.42	0.43
1:CA:321:A:O2'	1:CA:322:C:H5'	2.19	0.43
1:CA:652:U:H2'	1:CA:752:G:N1	2.33	0.43
1:CA:761:G:H2'	1:CA:762:C:H6	1.83	0.43
1:CA:865:A:O2'	1:CA:919:A:H5'	2.19	0.43
1:CA:961:U:C2'	1:CA:962:C:H5'	2.49	0.43
2:CB:24:TRP:CZ3	2:CB:26:PRO:HA	2.54	0.43
4:CD:15:GLU:CD	4:CD:15:GLU:N	2.72	0.43
4:CD:88:VAL:O	4:CD:88:VAL:HG12	2.19	0.43
6:CF:24:GLU:HG3	6:CF:25:ILE:HD13	2.00	0.43
6:CF:2:ARG:N	6:CF:67:MET:O	2.50	0.43
7:CG:23:VAL:C	7:CG:27:ILE:HD13	2.39	0.43
8:CH:88:LYS:O	8:CH:92:ARG:HD2	2.19	0.43
9:CI:118:LYS:CB	9:CI:118:LYS:NZ	2.81	0.43
15:CO:50:HIS:O	15:CO:53:HIS:CB	2.67	0.43
15:CO:39:LEU:HD12	15:CO:56:LEU:HD13	2.00	0.43
17:CQ:3:LYS:O	17:CQ:4:LYS:C	2.58	0.43
19:CS:43:GLU:O	19:CS:43:GLU:OE1	2.36	0.43
19:CS:45:VAL:C	19:CS:47:HIS:H	2.21	0.43
20:CT:12:ALA:C	20:CT:14:LYS:N	2.72	0.43
25:CY:160:GLU:HG3	25:CY:164:ILE:HD11	2.01	0.43
25:CY:31:GLY:O	25:CY:32:ARG:CB	2.67	0.43
26:D0:26:TYR:C	26:D0:67:VAL:HG11	2.39	0.43
28:D2:25:VAL:O	28:D2:28:LYS:N	2.52	0.43
34:D8:30:ARG:CG	34:D8:30:ARG:O	2.65	0.43
35:DA:818:G:H3'	35:DA:1187:G:H22	1.83	0.43
35:DA:1369:G:C2	35:DA:1370:C:C6	3.07	0.43
35:DA:1745(A):C:H5'	35:DA:1746:G:OP2	2.18	0.43
35:DA:1839:G:C4	35:DA:1840:G:C8	3.07	0.43
35:DA:1982:C:H2'	35:DA:1983:C:C6	2.54	0.43
35:DA:2075:U:C4	35:DA:2238:G:C6	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2684:U:C2'	35:DA:2685:G:H5'	2.49	0.43
35:DA:349:G:N3	35:DA:349:G:H2'	2.33	0.43
35:DA:382:G:H1	35:DA:392:C:H42	1.67	0.43
35:DA:28:A:C8	35:DA:513:A:C5	3.07	0.43
35:DA:71:A:H3'	35:DA:71:A:OP2	2.19	0.43
35:DA:751:A:H62	35:DA:789:A:H62	1.66	0.43
35:DA:863:A:OP1	47:DQ:21:THR:HB	2.19	0.43
38:DD:91:ARG:HG2	38:DD:91:ARG:NH1	2.33	0.43
38:DD:81:ALA:H	38:DD:94:LEU:CD1	2.32	0.43
39:DE:9:VAL:HG23	39:DE:9:VAL:O	2.18	0.43
41:DG:115:ARG:NH1	41:DG:136:ARG:HD2	2.33	0.43
41:DG:125:PHE:HA	41:DG:131:TYR:HB2	2.01	0.43
41:DG:16:ARG:HG3	41:DG:16:ARG:HH11	1.83	0.43
41:DG:40:ASN:CG	41:DG:41:GLN:H	2.23	0.43
42:DH:144:VAL:CG1	42:DH:144:VAL:O	2.65	0.43
43:DI:42:SER:C	43:DI:44:LEU:N	2.71	0.43
44:DN:17:ASP:OD2	44:DN:56:ASN:HB3	2.19	0.43
45:DO:108:GLU:O	45:DO:109:LYS:C	2.57	0.43
47:DQ:22:LYS:NZ	47:DQ:22:LYS:CA	2.82	0.43
35:DA:2469:A:O2'	47:DQ:56:ARG:HG3	2.19	0.43
48:DR:10:LEU:HD13	48:DR:17:ARG:CZ	2.49	0.43
49:DS:36:TYR:O	49:DS:37:ALA:HB2	2.19	0.43
49:DS:54:LEU:O	49:DS:56:LEU:N	2.51	0.43
49:DS:77:ALA:O	49:DS:80:LEU:N	2.44	0.43
49:DS:28:VAL:O	49:DS:89:ARG:HG2	2.19	0.43
51:DU:65:ILE:H	51:DU:65:ILE:CD1	2.31	0.43
51:DU:92:ARG:CG	51:DU:92:ARG:O	2.67	0.43
52:DV:27:ALA:O	52:DV:28:GLU:C	2.58	0.43
56:DZ:148:ASP:O	56:DZ:149:SER:CB	2.66	0.43
56:DZ:57:ILE:N	56:DZ:57:ILE:CD1	2.79	0.43
56:DZ:77:ASP:O	56:DZ:78:LYS:C	2.58	0.43
56:DZ:79:ARG:C	56:DZ:80:ARG:HG2	2.39	0.43
1:AA:1115:C:H2'	1:AA:1116:C:C6	2.54	0.42
1:AA:1459:C:C2'	1:AA:1460:A:H5'	2.49	0.42
1:AA:252:U:H5'	1:AA:253:U:OP2	2.19	0.42
1:AA:619:U:O2	4:AD:135:LEU:HD23	2.19	0.42
1:AA:651:C:O2'	1:AA:652:U:H5'	2.19	0.42
1:AA:797:C:H2'	1:AA:798:G:H8	1.84	0.42
1:AA:828:A:C8	1:AA:859:A:C4	3.07	0.42
1:AA:828:A:C8	1:AA:859:A:C5	3.07	0.42
2:AB:145:LEU:HD13	2:AB:145:LEU:C	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:112:VAL:HG22	2:AB:149:LEU:HD22	2.01	0.42
2:AB:185:ILE:HG22	2:AB:199:TYR:CB	2.45	0.42
3:AC:150:LYS:CB	3:AC:201:TYR:HB2	2.40	0.42
4:AD:56:VAL:C	4:AD:58:LEU:H	2.23	0.42
5:AE:76:ILE:CG2	5:AE:78:HIS:O	2.67	0.42
5:AE:91:LEU:HD23	5:AE:110:LEU:HD11	2.00	0.42
6:AF:82:ARG:HH11	6:AF:82:ARG:HA	1.84	0.42
7:AG:143:ARG:CB	7:AG:143:ARG:NH1	2.81	0.42
8:AH:119:LEU:HB2	8:AH:120:THR:H	1.72	0.42
9:AI:116:LYS:HA	9:AI:121:ARG:O	2.18	0.42
9:AI:92:TYR:HB3	9:AI:95:LYS:HD2	2.00	0.42
1:AA:1307:U:O4'	13:AM:109:THR:HG21	2.19	0.42
14:AN:39:LEU:HD11	14:AN:47:LEU:CD1	2.48	0.42
16:AP:12:LYS:O	16:AP:13:HIS:CB	2.66	0.42
16:AP:14:ASN:OD1	16:AP:16:HIS:HE1	2.02	0.42
16:AP:39:TYR:HB2	16:AP:49:LEU:HB2	1.99	0.42
18:AR:29:PHE:HD1	18:AR:39:VAL:HG11	1.83	0.42
18:AR:53:ARG:NH2	18:AR:60:ALA:CA	2.82	0.42
20:AT:78:ALA:O	20:AT:79:ARG:C	2.56	0.42
25:AY:143:LEU:C	25:AY:145:LYS:H	2.22	0.42
25:AY:149:LEU:HA	25:AY:153:GLU:OE1	2.19	0.42
25:AY:36:ALA:C	25:AY:38:LEU:H	2.23	0.42
25:AY:43:VAL:O	25:AY:50:VAL:HG22	2.19	0.42
28:B2:22:GLU:O	28:B2:25:VAL:HG13	2.18	0.42
29:B3:26:LEU:O	29:B3:28:LEU:HG	2.19	0.42
33:B7:16:HIS:CE1	35:BA:465:G:C4'	3.02	0.42
33:B7:29:LYS:NZ	33:B7:32:LYS:HZ2	2.17	0.42
35:BA:1495:A:OP1	35:BA:1495:A:O4'	2.37	0.42
35:BA:1553:A:H4'	35:BA:1553:A:OP1	2.19	0.42
35:BA:1271:G:N2	35:BA:1617:C:H4'	2.34	0.42
35:BA:195:A:N7	35:BA:197:A:OP1	2.52	0.42
35:BA:2223:G:H2'	35:BA:2224:G:C5'	2.48	0.42
35:BA:271(L):U:H5''	35:BA:271(M):G:C4	2.54	0.42
35:BA:2805:G:H2'	35:BA:2807:G:C8	2.53	0.42
35:BA:324:A:N6	35:BA:338:G:O2'	2.50	0.42
35:BA:491:G:C4	35:BA:492:A:C8	3.07	0.42
35:BA:567:A:N1	35:BA:568:U:O2	2.51	0.42
35:BA:670:A:H5'	35:BA:671:C:OP1	2.19	0.42
35:BA:811:U:OP2	46:BP:24:GLY:HA2	2.19	0.42
35:BA:915:C:H2'	35:BA:916:G:H8	1.83	0.42
35:BA:1566:A:OP1	38:BD:211:ARG:NE	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BD:260:ARG:CZ	38:BD:264:LYS:HD3	2.49	0.42
38:BD:34:VAL:CG2	38:BD:35:LYS:NZ	2.82	0.42
38:BD:82:ILE:HA	38:BD:92:ILE:O	2.19	0.42
35:BA:2572:A:C6	39:BE:144:ARG:NH2	2.87	0.42
39:BE:154:LYS:CE	39:BE:154:LYS:HA	2.45	0.42
40:BF:197:ASP:O	40:BF:200:GLU:HB3	2.19	0.42
40:BF:3:GLU:O	40:BF:19:GLU:CB	2.67	0.42
35:BA:660:G:C5'	40:BF:99:TYR:CE2	3.01	0.42
41:BG:102:PHE:O	41:BG:106:LEU:HB3	2.19	0.42
41:BG:166:ASP:OD1	41:BG:166:ASP:C	2.57	0.42
42:BH:151:ILE:HD13	42:BH:151:ILE:H	1.83	0.42
44:BN:57:ALA:H	44:BN:124:ALA:HA	1.84	0.42
45:BO:13:ASN:ND2	45:BO:97:ARG:HG2	2.34	0.42
45:BO:81:ASP:CG	45:BO:81:ASP:O	2.57	0.42
46:BP:108:LYS:C	46:BP:110:TYR:N	2.70	0.42
48:BR:20:LEU:CG	48:BR:21:TYR:N	2.81	0.42
48:BR:30:THR:OG1	48:BR:75:LEU:HD21	2.19	0.42
48:BR:73:VAL:HG23	48:BR:74:LYS:HD3	2.00	0.42
50:BT:28:VAL:HG21	50:BT:46:GLU:OE2	2.19	0.42
50:BT:87:ASP:OD2	50:BT:90:GLN:HG3	2.19	0.42
51:BU:35:ALA:O	51:BU:36:ARG:C	2.57	0.42
52:BV:39:LEU:O	52:BV:50:PRO:HA	2.19	0.42
53:BW:65:LEU:HD13	53:BW:68:ARG:HD2	2.01	0.42
54:BX:40:LYS:C	54:BX:42:ALA:N	2.71	0.42
54:BX:32:PRO:HB3	54:BX:72:LYS:HD2	2.01	0.42
54:BX:77:LYS:HA	54:BX:77:LYS:CE	2.46	0.42
1:CA:1050:G:H2'	1:CA:1051:C:C6	2.54	0.42
1:CA:1059:C:O2	10:CJ:53:PRO:HG3	2.19	0.42
1:CA:1088:G:C4	1:CA:1089:G:N7	2.87	0.42
1:CA:1135:U:H2'	1:CA:1137:C:O2	2.18	0.42
1:CA:113:G:H2'	1:CA:114:U:H6	1.84	0.42
1:CA:1210:C:H4'	1:CA:1214:C:N4	2.34	0.42
1:CA:960:U:C4	1:CA:1225:A:H1'	2.54	0.42
1:CA:66:G:O4'	1:CA:173:U:C4	2.72	0.42
1:CA:18:C:H2'	1:CA:19:C:H6	1.84	0.42
1:CA:27:G:H2'	1:CA:28:G:C8	2.53	0.42
1:CA:809:G:C2'	1:CA:810:C:O5'	2.66	0.42
1:CA:570:G:C5	1:CA:873:A:C6	3.07	0.42
2:CB:84:GLU:O	2:CB:219:VAL:HG11	2.19	0.42
2:CB:87:ARG:O	2:CB:223:ILE:HD11	2.19	0.42
4:CD:171:GLY:HA2	4:CD:172:PRO:HD3	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CD:19:LEU:O	4:CD:21:LEU:N	2.46	0.42
4:CD:33:MET:HE3	4:CD:33:MET:HA	2.01	0.42
4:CD:13:ARG:HH22	4:CD:36:ARG:HH11	1.66	0.42
6:CF:27:GLN:O	6:CF:31:GLU:HG3	2.19	0.42
6:CF:53:ALA:O	6:CF:54:LYS:HB2	2.18	0.42
8:CH:3:THR:O	8:CH:5:PRO:HD3	2.19	0.42
8:CH:78:GLN:OE1	8:CH:78:GLN:HA	2.19	0.42
10:CJ:47:PHE:HD1	10:CJ:47:PHE:O	2.01	0.42
14:CN:39:LEU:HD11	14:CN:47:LEU:CD1	2.49	0.42
17:CQ:67:LYS:C	17:CQ:70:ARG:HH12	2.22	0.42
25:CY:149:LEU:HA	25:CY:153:GLU:OE1	2.18	0.42
25:CY:62:ASP:OD1	25:CY:65:THR:HB	2.19	0.42
27:D1:56:GLN:O	27:D1:57:GLU:HG2	2.19	0.42
27:D1:76:ARG:C	27:D1:78:LYS:NZ	2.72	0.42
31:D5:40:LYS:NZ	31:D5:45:VAL:CA	2.76	0.42
35:DA:1050:A:O2'	35:DA:2752:C:H1'	2.19	0.42
35:DA:1368:G:C6	35:DA:1369:G:N7	2.87	0.42
35:DA:1605:C:C5'	35:DA:1610:A:N6	2.82	0.42
35:DA:1654:A:OP1	48:DR:2:ARG:HB2	2.19	0.42
35:DA:1806:C:H42	35:DA:1812:A:N6	2.17	0.42
35:DA:1997:G:C2	35:DA:1998:G:C5	3.07	0.42
35:DA:200:U:H2'	35:DA:201:C:C5'	2.47	0.42
35:DA:2080:G:H2'	35:DA:2081:C:C6	2.54	0.42
35:DA:19:C:H2'	35:DA:20:C:H6	1.84	0.42
35:DA:2223:G:O2'	35:DA:2224:G:H5'	2.19	0.42
35:DA:2262:U:C3'	35:DA:2263:C:H5''	2.49	0.42
35:DA:2283:C:C6	35:DA:2389:G:H2'	2.53	0.42
35:DA:2415:G:C2	35:DA:2416:C:C2	3.07	0.42
35:DA:2727:G:C5	35:DA:2728:U:H5	2.37	0.42
35:DA:2738:A:C2	35:DA:2739:U:C6	3.07	0.42
35:DA:2763:G:H5'	35:DA:2763:G:H8	1.83	0.42
35:DA:2832:U:C5	35:DA:2884:U:H5''	2.54	0.42
35:DA:325:G:H2'	35:DA:326:G:C8	2.50	0.42
35:DA:58:G:N3	35:DA:73:A:H2	2.17	0.42
35:DA:779:U:OP1	38:DD:49:ILE:HG23	2.19	0.42
35:DA:809:G:C4'	35:DA:1254:A:H1'	2.48	0.42
35:DA:870:A:N1	35:DA:871:U:C2	2.87	0.42
35:DA:898:C:N3	35:DA:899:A:H1'	2.34	0.42
35:DA:94(A):G:H2'	35:DA:95:G:O4'	2.19	0.42
38:DD:211:ARG:HA	38:DD:214:TRP:CE3	2.54	0.42
39:DE:3:GLY:CA	39:DE:81:ILE:HG21	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DE:9:VAL:HG11	39:DE:25:VAL:HG12	2.00	0.42
40:DF:114:VAL:CG2	40:DF:115:ALA:N	2.73	0.42
40:DF:164:ARG:O	40:DF:165:ARG:C	2.58	0.42
40:DF:17:ARG:HG3	40:DF:17:ARG:NH1	2.33	0.42
41:DG:138:GLN:HB3	41:DG:152:LEU:HD23	2.01	0.42
43:DI:132:PRO:O	43:DI:135:GLU:HG2	2.19	0.42
43:DI:51:ILE:HA	43:DI:51:ILE:HD13	1.75	0.42
43:DI:79:ILE:C	43:DI:81:VAL:H	2.21	0.42
44:DN:26:LEU:O	44:DN:30:ILE:HG13	2.19	0.42
44:DN:35:ARG:HD3	44:DN:37:LYS:HD3	2.00	0.42
46:DP:108:LYS:C	46:DP:110:TYR:N	2.71	0.42
46:DP:39:LYS:CD	46:DP:40:SER:N	2.65	0.42
48:DR:17:ARG:CG	48:DR:17:ARG:NH1	2.81	0.42
48:DR:71:GLN:HB2	48:DR:71:GLN:HE21	1.58	0.42
50:DT:74:ARG:HG2	50:DT:74:ARG:NH1	2.32	0.42
51:DU:31:SER:HB3	51:DU:34:LYS:CG	2.48	0.42
52:DV:51:VAL:HG12	52:DV:52:VAL:N	2.33	0.42
53:DW:44:ALA:O	53:DW:45:TYR:C	2.57	0.42
1:AA:1017:G:C2	1:AA:1018:C:N3	2.88	0.42
1:AA:320:C:HO2'	1:AA:1435:G:H1'	1.84	0.42
1:AA:1433:A:C8	1:AA:1468:A:C6	3.06	0.42
1:AA:264:U:O2'	17:AQ:64:PRO:HB2	2.18	0.42
1:AA:474:G:H2'	1:AA:475:G:H8	1.83	0.42
1:AA:536:C:H2'	1:AA:537:G:C8	2.53	0.42
1:AA:651:C:H2'	1:AA:652:U:H6	1.83	0.42
1:AA:68:G:N2	1:AA:69:G:H1'	2.33	0.42
1:AA:714:G:N2	1:AA:777:A:H1'	2.27	0.42
1:AA:828:A:N7	1:AA:859:A:C8	2.86	0.42
1:AA:91:C:H6	1:AA:91:C:O5'	2.01	0.42
2:AB:164:VAL:O	2:AB:165:VAL:O	2.36	0.42
2:AB:170:GLU:C	2:AB:172:ILE:N	2.71	0.42
2:AB:84:GLU:O	2:AB:219:VAL:HG11	2.19	0.42
2:AB:51:LEU:O	2:AB:55:PHE:CD2	2.72	0.42
2:AB:75:LYS:HE3	2:AB:75:LYS:CA	2.49	0.42
4:AD:3:ARG:HG2	4:AD:3:ARG:HH21	1.85	0.42
5:AE:129:ILE:HG22	5:AE:130:ASN:N	2.33	0.42
7:AG:143:ARG:HH11	7:AG:143:ARG:HB3	1.84	0.42
8:AH:14:ARG:NH1	8:AH:14:ARG:HB3	2.34	0.42
9:AI:3:GLN:HG2	9:AI:20:ARG:HE	1.84	0.42
10:AJ:94:VAL:CG1	10:AJ:95:GLU:N	2.81	0.42
12:AL:41:ARG:HH11	12:AL:41:ARG:CB	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:116:THR:HG22	13:AM:117:VAL:H	1.80	0.42
1:AA:1060:C:C5'	14:AN:45:ARG:HH22	2.15	0.42
17:AQ:17:LYS:HA	17:AQ:46:ASP:O	2.18	0.42
17:AQ:67:LYS:O	17:AQ:69:LYS:N	2.52	0.42
17:AQ:92:ARG:O	17:AQ:93:GLN:C	2.56	0.42
20:AT:12:ALA:C	20:AT:14:LYS:H	2.22	0.42
20:AT:24:LEU:O	20:AT:27:LYS:N	2.52	0.42
25:AY:129:ILE:O	25:AY:132:ILE:N	2.52	0.42
25:AY:52:LEU:C	25:AY:54:GLN:N	2.70	0.42
25:AY:76:LEU:HG	25:AY:99:LEU:HD21	2.01	0.42
27:B1:21:ARG:HH11	27:B1:21:ARG:CG	2.32	0.42
35:BA:1416:G:H1'	35:BA:1417:C:C6	2.54	0.42
35:BA:1613:G:H3'	35:BA:1614:A:H5''	2.01	0.42
35:BA:1669:A:H2'	35:BA:1670:C:H5'	2.01	0.42
35:BA:2009:G:N1	35:BA:2010:G:N7	2.67	0.42
35:BA:2505:G:H2'	35:BA:2576:G:C6	2.54	0.42
35:BA:2539:C:O2	35:BA:2539:C:H2'	2.20	0.42
35:BA:704:G:N3	35:BA:726:G:C2	2.87	0.42
35:BA:744:G:O2'	35:BA:745:G:H5'	2.19	0.42
35:BA:982:C:H6	35:BA:982:C:O5'	2.02	0.42
37:BC:44:HIS:CD2	37:BC:175:VAL:HA	2.54	0.42
38:BD:105:ILE:O	38:BD:105:ILE:HG23	2.19	0.42
38:BD:13:ARG:HA	38:BD:13:ARG:HD2	1.95	0.42
38:BD:164:GLN:C	38:BD:165:ILE:HD12	2.40	0.42
38:BD:45:ASN:OD1	38:BD:48:ARG:O	2.37	0.42
38:BD:72:LYS:HB2	38:BD:97:TYR:HE2	1.84	0.42
38:BD:88:ARG:HH11	38:BD:88:ARG:HG2	1.84	0.42
39:BE:28:ALA:O	39:BE:180:ASN:OD1	2.37	0.42
40:BF:109:GLY:C	40:BF:111:ALA:N	2.72	0.42
40:BF:175:THR:HG23	40:BF:175:THR:O	2.19	0.42
35:BA:615:G:OP2	40:BF:43:LYS:HE2	2.20	0.42
41:BG:101:ILE:HD11	41:BG:105:LYS:CE	2.43	0.42
41:BG:14:GLU:O	41:BG:18:GLU:CB	2.66	0.42
42:BH:46:GLU:CG	42:BH:51:ARG:HB2	2.50	0.42
44:BN:15:LEU:HD12	44:BN:136:GLU:CB	2.48	0.42
50:BT:51:ARG:O	50:BT:52:ILE:HG23	2.19	0.42
44:BN:41:ASP:CA	51:BU:64:ARG:NH1	2.82	0.42
52:BV:61:VAL:CG1	52:BV:62:LEU:N	2.79	0.42
54:BX:37:THR:HG23	54:BX:54:VAL:HB	2.00	0.42
55:BY:31:LEU:HD12	55:BY:34:LYS:N	2.35	0.42
55:BY:43:ASN:CG	55:BY:64:GLU:HA	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:BZ:139:VAL:HG12	56:BZ:141:VAL:N	2.34	0.42
56:BZ:44:PHE:C	56:BZ:44:PHE:CD1	2.93	0.42
1:CA:1302:U:O2'	1:CA:1303:C:H5'	2.19	0.42
1:CA:1316:G:H2'	1:CA:1317:C:H5'	2.01	0.42
1:CA:514:C:N4	1:CA:537:G:H1	2.15	0.42
1:CA:553:A:O2'	1:CA:554:C:H5'	2.20	0.42
1:CA:61:G:H2'	1:CA:62:U:O4'	2.19	0.42
1:CA:778:G:C2'	1:CA:779:C:H5'	2.49	0.42
1:CA:939:G:H2'	1:CA:940:C:C6	2.52	0.42
2:CB:100:GLY:O	2:CB:104:ASN:C	2.58	0.42
2:CB:105:PHE:O	2:CB:106:LYS:C	2.57	0.42
2:CB:213:LEU:HD22	2:CB:214:ILE:HG12	2.01	0.42
4:CD:152:SER:C	4:CD:154:ASN:H	2.22	0.42
4:CD:177:ASP:O	4:CD:177:ASP:OD1	2.37	0.42
5:CE:144:THR:C	5:CE:148:VAL:HG23	2.38	0.42
7:CG:143:ARG:HH11	7:CG:143:ARG:HB3	1.82	0.42
7:CG:143:ARG:CB	7:CG:143:ARG:NH1	2.78	0.42
7:CG:63:LYS:HD2	7:CG:63:LYS:HA	1.85	0.42
8:CH:7:ALA:HA	8:CH:10:LEU:HG	2.00	0.42
9:CI:8:GLY:HA3	9:CI:15:ALA:HB3	2.01	0.42
9:CI:53:VAL:HG12	9:CI:95:LYS:CE	2.49	0.42
9:CI:83:ARG:C	9:CI:86:VAL:HG12	2.39	0.42
10:CJ:98:ILE:HG22	10:CJ:99:LYS:N	2.34	0.42
14:CN:33:VAL:HG12	14:CN:34:TYR:H	1.84	0.42
15:CO:43:LEU:C	15:CO:45:VAL:N	2.72	0.42
16:CP:22:THR:HA	16:CP:33:ILE:HG12	2.01	0.42
17:CQ:10:VAL:HG11	17:CQ:53:LEU:HA	2.00	0.42
6:CF:100:ASN:O	18:CR:28:GLU:HB3	2.19	0.42
20:CT:16:HIS:O	20:CT:19:SER:N	2.52	0.42
20:CT:84:LEU:HD13	20:CT:84:LEU:C	2.38	0.42
25:CY:130:ARG:HG2	25:CY:130:ARG:NH1	2.35	0.42
29:D3:58:VAL:HG12	29:D3:59:VAL:N	2.33	0.42
35:DA:1208:C:C5	35:DA:1209:G:N7	2.87	0.42
35:DA:1210:A:H1'	35:DA:1212:G:C4	2.54	0.42
35:DA:1790:C:H3'	35:DA:1828:G:H22	1.84	0.42
35:DA:2052:G:C4	35:DA:2053:G:C8	3.07	0.42
35:DA:919:G:N2	35:DA:2269:A:OP2	2.47	0.42
35:DA:2464:C:O2'	35:DA:2465:C:P	2.76	0.42
35:DA:2473:U:O2	35:DA:2473:U:H2'	2.19	0.42
35:DA:258:G:C6	35:DA:259:G:C5	3.07	0.42
35:DA:2618:G:O2'	35:DA:2619:C:H5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2685:G:OP1	50:DT:51:ARG:NH2	2.51	0.42
35:DA:2716:U:O2'	35:DA:2717:G:H5'	2.19	0.42
35:DA:2723:C:HO2'	35:DA:2724:C:H5'	1.84	0.42
35:DA:2851:A:C6	35:DA:2852:G:C6	3.07	0.42
35:DA:460:A:C2	35:DA:470:A:C4	3.07	0.42
35:DA:629:G:O2'	35:DA:630:G:H5'	2.19	0.42
35:DA:776:G:O6	35:DA:793:A:H2'	2.19	0.42
35:DA:817:C:H4'	35:DA:932:G:C6	2.54	0.42
36:DB:82:G:H2'	36:DB:83:G:C8	2.50	0.42
38:DD:34:VAL:CG2	38:DD:35:LYS:NZ	2.82	0.42
38:DD:77:ALA:CB	38:DD:97:TYR:CD1	3.02	0.42
40:DF:139:PHE:O	40:DF:142:TRP:HB3	2.18	0.42
40:DF:46:ARG:HG3	40:DF:48:THR:CG2	2.47	0.42
41:DG:101:ILE:HG23	41:DG:102:PHE:H	1.84	0.42
41:DG:38:VAL:CG1	41:DG:91:ARG:HD3	2.48	0.42
42:DH:127:GLU:HB2	42:DH:130:ARG:HB3	2.01	0.42
42:DH:79:VAL:HG23	42:DH:80:SER:N	2.34	0.42
43:DI:48:GLU:CD	43:DI:48:GLU:C	2.77	0.42
44:DN:36:GLY:HA3	44:DN:48:MET:SD	2.60	0.42
44:DN:82:LEU:H	44:DN:82:LEU:HD12	1.83	0.42
46:DP:29:LYS:N	46:DP:29:LYS:CD	2.82	0.42
47:DQ:38:GLU:HB2	47:DQ:127:ILE:HG23	2.01	0.42
50:DT:100:TYR:O	50:DT:101:PHE:C	2.57	0.42
51:DU:17:ILE:CA	51:DU:20:LEU:HD23	2.37	0.42
51:DU:46:ALA:O	51:DU:49:HIS:HB2	2.18	0.42
51:DU:64:ARG:CA	51:DU:64:ARG:NH2	2.82	0.42
51:DU:91:ASP:OD2	51:DU:96:ALA:CB	2.67	0.42
53:DW:20:VAL:O	53:DW:21:VAL:C	2.58	0.42
54:DX:83:VAL:C	54:DX:85:PRO:HD3	2.40	0.42
56:DZ:8:TYR:CA	56:DZ:62:PRO:HG2	2.47	0.42
1:AA:1148:U:H2'	1:AA:1149:C:O4'	2.20	0.42
1:AA:1220:G:H21	19:AS:54:GLY:HA2	1.84	0.42
1:AA:125:U:H2'	1:AA:126:G:H8	1.84	0.42
1:AA:1405:G:O2'	1:AA:1406:U:C5'	2.67	0.42
1:AA:385:C:H2'	1:AA:386:C:C6	2.53	0.42
1:AA:777:A:H2'	1:AA:778:G:H8	1.84	0.42
1:AA:696:A:H1'	1:AA:786:G:O2'	2.19	0.42
4:AD:101:LEU:O	4:AD:104:VAL:N	2.52	0.42
4:AD:138:TYR:C	4:AD:138:TYR:HD2	2.23	0.42
4:AD:72:GLU:O	4:AD:76:ARG:N	2.48	0.42
6:AF:42:GLU:C	6:AF:44:GLY:H	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:26:PHE:O	7:AG:30:ILE:HG12	2.20	0.42
10:AJ:22:LYS:C	10:AJ:22:LYS:HD2	2.39	0.42
10:AJ:48:THR:CB	10:AJ:62:HIS:HB3	2.49	0.42
11:AK:116:HIS:O	11:AK:117:ASN:HB3	2.19	0.42
23:AW:10:G:H22	23:AW:27:G:C1'	2.25	0.42
25:AY:52:LEU:HD23	25:AY:52:LEU:O	2.19	0.42
26:B0:23:VAL:CG1	26:B0:24:LYS:N	2.82	0.42
29:B3:22:ALA:HB1	29:B3:46:ASN:HD22	1.84	0.42
35:BA:139(A):G:N2	54:BX:44:GLU:OE1	2.52	0.42
35:BA:1666:G:C2'	35:BA:1667:G:H5'	2.49	0.42
35:BA:1906:G:O2'	35:BA:1907:G:H5'	2.19	0.42
35:BA:203:C:C6	35:BA:204:A:C8	3.07	0.42
35:BA:19:C:H2'	35:BA:20:C:H6	1.83	0.42
35:BA:2114:A:C3'	35:BA:2115:G:H5'	2.49	0.42
35:BA:2208:A:H1'	35:BA:2219:G:N3	2.34	0.42
35:BA:2261:C:O4'	35:BA:2388:A:H1'	2.19	0.42
35:BA:229:A:H3'	35:BA:230:U:C5'	2.42	0.42
35:BA:2285:C:H42	35:BA:2383:G:H1	1.67	0.42
35:BA:2483:C:N3	47:BQ:124:LYS:NZ	2.65	0.42
35:BA:2545:G:C2	35:BA:2546:U:C2	3.06	0.42
35:BA:2571:C:O2	35:BA:2571:C:C2'	2.67	0.42
35:BA:2574:G:C5	35:BA:2575:C:C5	3.07	0.42
35:BA:2704:C:C2	35:BA:2705:A:C8	3.07	0.42
35:BA:372:G:HO2'	35:BA:373:U:P	2.41	0.42
35:BA:593:G:H2'	35:BA:594:U:H6	1.83	0.42
35:BA:64:A:C2	35:BA:65:C:C2	3.08	0.42
35:BA:675:A:C6	35:BA:676:A:C6	3.07	0.42
35:BA:783:A:H2'	35:BA:784:A:C4'	2.47	0.42
35:BA:937:U:H2'	35:BA:938:G:O4'	2.19	0.42
36:BB:115:G:H2'	36:BB:116:G:C8	2.54	0.42
36:BB:66:A:O2'	36:BB:67:G:P	2.77	0.42
37:BC:56:GLN:NE2	37:BC:169:GLY:H	2.17	0.42
38:BD:211:ARG:HA	38:BD:214:TRP:CE3	2.53	0.42
38:BD:265:PRO:CG	38:BD:266:SER:N	2.82	0.42
38:BD:79:VAL:O	38:BD:79:VAL:HG12	2.19	0.42
38:BD:94:LEU:O	38:BD:94:LEU:CD1	2.65	0.42
39:BE:201:THR:HG22	39:BE:202:LYS:N	2.34	0.42
39:BE:63:LEU:O	39:BE:65:GLY:N	2.52	0.42
40:BF:111:ALA:O	40:BF:112:MET:C	2.58	0.42
40:BF:160:ASN:ND2	40:BF:160:ASN:C	2.71	0.42
40:BF:178:PRO:O	40:BF:180:GLY:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BG:160:VAL:CG1	41:BG:161:THR:N	2.81	0.42
41:BG:4:ASP:O	41:BG:4:ASP:OD1	2.37	0.42
36:BB:43:C:H4'	41:BG:66:GLN:NE2	2.34	0.42
42:BH:84:SER:O	42:BH:133:VAL:O	2.37	0.42
43:BI:133:HIS:CG	43:BI:134:PRO:HD2	2.53	0.42
43:BI:25:TYR:O	43:BI:30:LEU:HG	2.20	0.42
43:BI:90:GLY:O	43:BI:91:SER:C	2.57	0.42
44:BN:110:GLY:O	44:BN:111:PRO:C	2.56	0.42
44:BN:3:THR:C	44:BN:5:VAL:N	2.71	0.42
44:BN:46:VAL:HG11	44:BN:48:MET:CG	2.47	0.42
44:BN:46:VAL:O	44:BN:47:ALA:CB	2.67	0.42
46:BP:82:GLY:HA2	46:BP:113:LYS:O	2.19	0.42
46:BP:34:GLY:O	46:BP:36:LYS:HG3	2.19	0.42
47:BQ:23:GLY:CA	47:BQ:101:ARG:HB2	2.48	0.42
47:BQ:70:PRO:HA	47:BQ:95:ALA:HB2	2.00	0.42
47:BQ:43:THR:CA	47:BQ:94:VAL:HG12	2.43	0.42
47:BQ:20:ALA:CB	47:BQ:99:PRO:HG2	2.37	0.42
49:BS:20:ARG:HA	49:BS:20:ARG:HD3	1.66	0.42
51:BU:66:ASN:OD1	51:BU:76:TYR:HB2	2.19	0.42
55:BY:81:LYS:HD3	55:BY:97:ARG:C	2.40	0.42
55:BY:76:CYS:CB	55:BY:96:ILE:HD11	2.49	0.42
56:BZ:163:LEU:HD23	56:BZ:163:LEU:H	1.81	0.42
56:BZ:56:VAL:HA	56:BZ:70:LEU:CG	2.49	0.42
56:BZ:61:LEU:HB2	56:BZ:65:GLN:C	2.39	0.42
1:CA:1197:G:O2'	1:CA:1198:G:H5'	2.19	0.42
1:CA:1348:U:H2'	1:CA:1349:A:H8	1.83	0.42
1:CA:1372:U:OP1	9:CI:71:SER:HB3	2.19	0.42
1:CA:1376:U:O2'	1:CA:1377:A:H5'	2.18	0.42
1:CA:353:A:H5'	1:CA:353:A:H8	1.84	0.42
1:CA:482:A:H3'	1:CA:483:C:C6	2.54	0.42
1:CA:974:A:C8	14:CN:31:ARG:NE	2.87	0.42
2:CB:187:LEU:HD22	2:CB:188:ALA:N	2.35	0.42
2:CB:30:ARG:HG3	2:CB:31:TYR:CD2	2.55	0.42
4:CD:161:ASN:O	4:CD:165:MET:HB2	2.19	0.42
4:CD:17:VAL:CG1	4:CD:18:LYS:N	2.79	0.42
5:CE:112:LEU:O	5:CE:114:GLY:N	2.52	0.42
6:CF:97:PHE:N	18:CR:30:ASP:OD1	2.52	0.42
8:CH:95:VAL:HG22	8:CH:131:GLY:O	2.20	0.42
9:CI:17:VAL:HG22	9:CI:63:ILE:CG2	2.46	0.42
9:CI:70:LYS:O	9:CI:73:GLN:HB2	2.19	0.42
9:CI:92:TYR:N	9:CI:92:TYR:CD1	2.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CL:48:PRO:HD2	12:CL:49:ASN:OD1	2.18	0.42
1:CA:522:C:N4	12:CL:53:ARG:NH2	2.59	0.42
13:CM:4:ILE:HG22	13:CM:5:ALA:H	1.84	0.42
15:CO:45:VAL:HG12	15:CO:46:HIS:ND1	2.34	0.42
15:CO:55:GLY:O	15:CO:59:MET:HG3	2.19	0.42
16:CP:22:THR:HG22	16:CP:32:TYR:CA	2.47	0.42
19:CS:33:THR:HG22	19:CS:49:ILE:HG22	2.01	0.42
21:CU:17:THR:O	21:CU:22:ARG:NH1	2.51	0.42
25:CY:86:SER:O	25:CY:88:LEU:N	2.51	0.42
27:D1:53:VAL:HG12	27:D1:58:ILE:HG22	2.00	0.42
27:D1:89:GLU:O	27:D1:93:GLU:OE1	2.36	0.42
28:D2:34:GLU:O	28:D2:36:ARG:N	2.51	0.42
35:DA:1277:G:H2'	35:DA:1278:A:O4'	2.19	0.42
35:DA:1526:G:H2'	35:DA:1527:G:C8	2.54	0.42
35:DA:1560:G:H2'	35:DA:1561:G:C8	2.54	0.42
35:DA:1664:A:OP1	35:DA:1665:A:OP2	2.36	0.42
35:DA:1805:U:C2	35:DA:1813:G:C2	3.08	0.42
35:DA:1797:C:N3	35:DA:1823:G:C6	2.87	0.42
35:DA:2023:G:C2	35:DA:2024:G:C5	3.06	0.42
35:DA:2228:G:C5	35:DA:2229:C:C5	3.07	0.42
35:DA:2374:C:O2'	35:DA:2375:G:H5'	2.19	0.42
35:DA:2408:U:C6	35:DA:2408:U:OP2	2.73	0.42
35:DA:2517:C:C6	35:DA:2542:A:N1	2.87	0.42
35:DA:2594:C:H2'	35:DA:2595:G:O4'	2.18	0.42
35:DA:2639:A:H3'	35:DA:2640:G:C5'	2.49	0.42
35:DA:2697:G:C2	35:DA:2711:A:C2	3.07	0.42
35:DA:2768:C:C2'	35:DA:2769:C:H5'	2.49	0.42
35:DA:2785:C:H2'	35:DA:2786:U:H6	1.83	0.42
35:DA:296:C:C2'	35:DA:297:C:H5'	2.50	0.42
35:DA:29:U:H2'	35:DA:30:G:H8	1.84	0.42
35:DA:648:G:O4'	35:DA:2351:G:H5''	2.19	0.42
35:DA:798:G:C6	35:DA:799:G:C6	3.07	0.42
35:DA:817:C:O2'	35:DA:839:U:OP1	2.36	0.42
35:DA:868:U:C4	35:DA:869:G:N7	2.88	0.42
35:DA:880:G:N1	35:DA:898:C:N4	2.68	0.42
38:DD:203:ASN:O	38:DD:204:ILE:C	2.57	0.42
38:DD:81:ALA:O	38:DD:94:LEU:HD12	2.19	0.42
39:DE:6:GLY:CA	39:DE:27:LEU:O	2.67	0.42
40:DF:74:ARG:H	40:DF:74:ARG:HD2	1.84	0.42
41:DG:80:PHE:O	41:DG:81:LYS:CB	2.68	0.42
42:DH:94:TYR:CE1	42:DH:160:LYS:HD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DH:76:VAL:O	42:DH:77:LYS:C	2.57	0.42
44:DN:15:LEU:HD13	44:DN:15:LEU:C	2.40	0.42
44:DN:46:VAL:O	44:DN:47:ALA:CB	2.65	0.42
44:DN:55:VAL:HG11	44:DN:127:ASP:H	1.84	0.42
46:DP:146:VAL:CG2	46:DP:147:LEU:H	2.04	0.42
46:DP:16:ARG:NH1	46:DP:18:ARG:CG	2.80	0.42
49:DS:90:GLY:HA2	49:DS:92:TYR:CG	2.53	0.42
50:DT:70:VAL:HG12	50:DT:71:GLY:O	2.18	0.42
52:DV:1:MET:HE2	52:DV:44:LYS:HB3	2.01	0.42
53:DW:65:LEU:HD13	53:DW:68:ARG:HD2	2.01	0.42
1:AA:1442(A):G:C2	50:BT:118:ARG:HB2	2.55	0.42
1:AA:189(I):G:H2'	1:AA:189(J):G:H8	1.83	0.42
1:AA:243:A:C2	1:AA:246:A:C8	3.06	0.42
1:AA:474:G:O2'	1:AA:475:G:H5'	2.20	0.42
1:AA:594:G:H2'	1:AA:595:G:O4'	2.20	0.42
1:AA:723:U:H5''	1:AA:724:G:OP2	2.18	0.42
1:AA:577:G:H1'	1:AA:816:A:N3	2.34	0.42
2:AB:113:HIS:C	2:AB:115:LEU:N	2.73	0.42
3:AC:33:LEU:O	3:AC:36:ASP:HB3	2.19	0.42
5:AE:135:THR:O	5:AE:138:ALA:HB3	2.19	0.42
5:AE:60:TYR:HE2	5:AE:64:ARG:NE	2.16	0.42
6:AF:96:PRO:HB3	18:AR:30:ASP:CG	2.39	0.42
7:AG:130:GLY:C	7:AG:132:GLY:H	2.22	0.42
7:AG:18:TYR:HE1	7:AG:44:TYR:CE2	2.38	0.42
7:AG:27:ILE:HG23	7:AG:40:ALA:N	2.34	0.42
7:AG:47:CYS:C	7:AG:58:PRO:HG3	2.40	0.42
12:AL:102:ARG:CD	12:AL:108:ALA:O	2.67	0.42
1:AA:537:G:OP1	12:AL:113:ARG:NH2	2.52	0.42
12:AL:125:PRO:O	12:AL:125:PRO:HG2	2.19	0.42
1:AA:522:C:N4	12:AL:53:ARG:NH2	2.58	0.42
13:AM:69:GLU:HB2	13:AM:72:ALA:HB3	2.01	0.42
23:AW:28:U:O4	23:AW:44:A:N1	2.53	0.42
25:AY:7:TYR:CD2	25:AY:160:GLU:HG2	2.54	0.42
29:B3:40:THR:HA	29:B3:44:ARG:CZ	2.48	0.42
32:B6:22:ALA:HB2	32:B6:39:TYR:CE2	2.54	0.42
35:BA:1003:G:N2	35:BA:1153:C:C2	2.87	0.42
35:BA:1019:U:H3	35:BA:1142(A):A:N6	2.03	0.42
35:BA:1153:C:C4	35:BA:1154:G:C2	3.07	0.42
35:BA:1340:U:C5	35:BA:1603:A:C1'	3.02	0.42
35:BA:1409:C:O2	35:BA:1594:G:N2	2.52	0.42
35:BA:1416:G:HO2'	35:BA:1417:C:H6	1.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1797:C:H2'	35:BA:1798:U:H6	1.84	0.42
35:BA:1952:A:C6	35:BA:1953:A:C6	3.07	0.42
35:BA:1956:U:C2'	35:BA:1957:C:H5'	2.49	0.42
35:BA:1982:C:H3'	35:BA:1982:C:OP1	2.20	0.42
35:BA:2066:C:O2'	35:BA:2067:G:H5'	2.18	0.42
35:BA:2256:G:H2'	35:BA:2257:U:C6	2.55	0.42
26:B0:16:SER:HB3	35:BA:2261:C:H3'	2.02	0.42
35:BA:2508:G:H2'	35:BA:2509:G:H8	1.83	0.42
35:BA:2658:C:N4	35:BA:2664:G:N2	2.68	0.42
35:BA:271(V):G:H2'	35:BA:271(W):G:C8	2.54	0.42
35:BA:2768:C:C2'	35:BA:2769:C:H5'	2.50	0.42
35:BA:404:C:C3'	35:BA:405:U:H5'	2.47	0.42
35:BA:479:A:H61	35:BA:503:A:H61	1.67	0.42
35:BA:481:G:O2'	35:BA:482:A:H8	2.02	0.42
35:BA:510:C:OP1	35:BA:511:U:OP2	2.37	0.42
35:BA:562:U:O2'	35:BA:563:G:H5''	2.19	0.42
34:B8:18:ALA:HB3	35:BA:651:G:H4'	2.01	0.42
35:BA:68:G:N3	35:BA:68:G:H2'	2.34	0.42
35:BA:704:G:O2'	35:BA:726:G:N2	2.46	0.42
35:BA:742:G:H1	35:BA:756:C:N4	2.17	0.42
36:BB:87:G:C2'	36:BB:88:C:H5''	2.49	0.42
36:BB:97:G:C2	36:BB:98:G:C8	3.07	0.42
38:BD:69:ARG:NH2	38:BD:128:GLY:O	2.53	0.42
38:BD:92:ILE:C	38:BD:92:ILE:CD1	2.88	0.42
41:BG:156:ASP:O	41:BG:157:ILE:C	2.57	0.42
41:BG:133:LEU:O	41:BG:157:ILE:O	2.36	0.42
41:BG:167:GLU:O	41:BG:170:ARG:HB3	2.19	0.42
42:BH:121:ILE:HG22	42:BH:133:VAL:CG1	2.49	0.42
43:BI:48:GLU:CD	43:BI:48:GLU:C	2.78	0.42
45:BO:6:THR:CG2	45:BO:7:TYR:N	2.74	0.42
45:BO:87:ILE:HG13	45:BO:91:LEU:HD13	2.00	0.42
46:BP:126:VAL:HG22	46:BP:145:PRO:CB	2.48	0.42
46:BP:29:LYS:CD	46:BP:29:LYS:N	2.82	0.42
48:BR:53:HIS:O	48:BR:56:LYS:HB3	2.19	0.42
50:BT:108:ARG:HH11	50:BT:108:ARG:CB	2.32	0.42
50:BT:51:ARG:O	50:BT:61:PHE:HB2	2.20	0.42
51:BU:10:ARG:O	51:BU:11:ARG:C	2.58	0.42
51:BU:95:LEU:HD23	51:BU:95:LEU:HA	1.89	0.42
51:BU:97:ASP:O	51:BU:100:VAL:HG23	2.19	0.42
52:BV:34:GLU:O	52:BV:62:LEU:HG	2.19	0.42
53:BW:50:VAL:HG13	53:BW:51:LEU:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:BW:65:LEU:HD22	53:BW:68:ARG:H	1.85	0.42
54:BX:60:ARG:HE	54:BX:74:PRO:CG	2.33	0.42
56:BZ:100:VAL:O	56:BZ:123:ASP:HA	2.19	0.42
1:CA:1006:C:H2'	1:CA:1007:C:H6	1.84	0.42
1:CA:106:C:O2'	1:CA:107:G:H5'	2.19	0.42
1:CA:1079:G:H2'	1:CA:1080:A:H8	1.82	0.42
1:CA:1157:A:H1'	1:CA:1181:G:H21	1.85	0.42
1:CA:1271:G:C5'	1:CA:1314:C:H5''	2.43	0.42
1:CA:1423:G:H5'	45:DO:49:ARG:NH2	2.34	0.42
1:CA:15:G:C4	1:CA:16:A:C8	3.08	0.42
1:CA:453:A:C6	1:CA:454:C:C4	3.08	0.42
1:CA:536:C:H2'	1:CA:537:G:C8	2.54	0.42
1:CA:600:C:C4	1:CA:639:G:N1	2.88	0.42
1:CA:659:U:C2	1:CA:660:G:C8	3.07	0.42
1:CA:807:A:H2'	1:CA:808:C:C6	2.54	0.42
1:CA:908:A:N3	1:CA:909:A:C8	2.87	0.42
2:CB:118:LEU:O	2:CB:119:GLU:C	2.58	0.42
2:CB:167:PRO:O	2:CB:170:GLU:N	2.42	0.42
4:CD:111:ALA:HA	4:CD:116:GLN:OE1	2.19	0.42
4:CD:120:LEU:CD1	4:CD:120:LEU:N	2.69	0.42
4:CD:150:GLU:O	4:CD:153:ARG:N	2.51	0.42
7:CG:97:GLN:O	7:CG:98:SER:C	2.57	0.42
9:CI:4:TYR:H	9:CI:4:TYR:HD1	1.67	0.42
11:CK:104:GLN:O	11:CK:106:LYS:N	2.52	0.42
15:CO:36:ILE:O	15:CO:37:ASN:C	2.57	0.42
19:CS:78:ARG:HH11	19:CS:78:ARG:HG3	1.84	0.42
25:CY:76:LEU:O	25:CY:79:ILE:HB	2.20	0.42
27:D1:26:ARG:HB3	27:D1:34:THR:OG1	2.20	0.42
27:D1:30:VAL:O	27:D1:30:VAL:HG12	2.19	0.42
27:D1:56:GLN:NE2	27:D1:57:GLU:HB2	2.34	0.42
29:D3:1:MET:O	29:D3:3:ARG:N	2.53	0.42
32:D6:15:GLU:CD	32:D6:18:ARG:HG3	2.40	0.42
35:DA:83:G:C4	35:DA:102:G:N2	2.88	0.42
35:DA:1187:G:H8	35:DA:1187:G:O5'	2.03	0.42
35:DA:1283:G:N2	35:DA:1285:G:H3'	2.34	0.42
35:DA:13:A:N6	35:DA:525:U:C5	2.87	0.42
35:DA:1495:A:OP1	35:DA:1495:A:O4'	2.37	0.42
35:DA:1710:C:H2'	35:DA:1711:C:H6	1.82	0.42
35:DA:1847:A:N3	35:DA:1847:A:C2'	2.81	0.42
35:DA:1853:A:H2'	35:DA:1854:A:C8	2.54	0.42
35:DA:1973:G:H2'	35:DA:1974:C:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:201:C:H2'	35:DA:202:U:H5'	2.00	0.42
35:DA:213:A:C2'	35:DA:214:G:H5'	2.50	0.42
35:DA:2548:G:C6	35:DA:2561:A:N1	2.87	0.42
35:DA:2704:C:C5	35:DA:2705:A:N7	2.87	0.42
35:DA:39:C:C4	35:DA:40:C:N4	2.87	0.42
35:DA:537:C:C2	35:DA:538:G:C8	3.06	0.42
35:DA:618:C:H2'	35:DA:619:G:C8	2.54	0.42
35:DA:737:C:H2'	35:DA:738:G:O5'	2.18	0.42
35:DA:749:C:C4	35:DA:1618:A:C2	3.07	0.42
35:DA:920:G:O2'	35:DA:921:G:H5'	2.19	0.42
36:DB:70:C:O2'	36:DB:71:C:H5'	2.17	0.42
39:DE:44:TYR:CE2	39:DE:46:ALA:HB2	2.55	0.42
39:DE:53:PRO:O	39:DE:55:ASN:OD1	2.36	0.42
40:DF:83:PHE:O	40:DF:84:VAL:C	2.57	0.42
41:DG:102:PHE:HD1	41:DG:106:LEU:HD22	1.83	0.42
41:DG:18:GLU:O	41:DG:22:ARG:HB3	2.19	0.42
41:DG:70:VAL:HG23	41:DG:70:VAL:O	2.20	0.42
42:DH:94:TYR:CZ	42:DH:160:LYS:HD2	2.55	0.42
43:DI:38:LEU:HD12	43:DI:38:LEU:N	2.11	0.42
44:DN:70:LYS:HG2	44:DN:87:LEU:HD23	2.02	0.42
45:DO:102:VAL:HB	45:DO:106:LEU:HD11	2.02	0.42
45:DO:6:THR:O	45:DO:20:MET:HA	2.19	0.42
46:DP:35:HIS:O	46:DP:35:HIS:CD2	2.72	0.42
47:DQ:70:PRO:O	47:DQ:71:ASP:HB3	2.19	0.42
49:DS:74:ALA:HB2	49:DS:101:LEU:HD11	2.00	0.42
50:DT:28:VAL:HG11	50:DT:46:GLU:CD	2.38	0.42
50:DT:80:SER:O	50:DT:81:PRO:C	2.55	0.42
51:DU:57:PHE:CD1	51:DU:60:LEU:HD12	2.54	0.42
44:DN:41:ASP:CA	51:DU:64:ARG:NH1	2.83	0.42
52:DV:20:LEU:N	52:DV:20:LEU:HD12	2.35	0.42
52:DV:36:PRO:HG2	52:DV:60:GLU:CD	2.39	0.42
55:DY:31:LEU:CD1	55:DY:34:LYS:N	2.79	0.42
56:DZ:40:ASP:OD1	56:DZ:42:VAL:HG13	2.18	0.42
56:DZ:56:VAL:CG1	56:DZ:57:ILE:N	2.82	0.42
1:AA:1155:G:H2'	1:AA:1156:G:H5'	2.01	0.42
1:AA:1271:G:C5'	1:AA:1314:C:H5'	2.42	0.42
1:AA:27:G:H2'	1:AA:28:G:C8	2.54	0.42
1:AA:39:G:O2'	1:AA:40:C:H5'	2.20	0.42
1:AA:774:G:H5'	38:BD:202:LYS:NZ	2.34	0.42
1:AA:896:C:O2'	1:AA:897:C:H5'	2.19	0.42
2:AB:36:ARG:HG3	2:AB:37:ASN:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:51:LEU:HB3	2:AB:55:PHE:CE2	2.51	0.42
3:AC:58:GLU:HB2	3:AC:65:ALA:HB2	2.00	0.42
4:AD:192:GLU:C	4:AD:194:LEU:N	2.72	0.42
4:AD:30:LYS:HB3	4:AD:35:ARG:HD2	2.01	0.42
5:AE:67:VAL:HG22	5:AE:68:GLU:O	2.20	0.42
7:AG:133:GLY:HA2	7:AG:136:LYS:HG2	2.01	0.42
5:AE:152:ARG:HG2	8:AH:43:GLY:O	2.18	0.42
8:AH:78:GLN:HA	8:AH:78:GLN:OE1	2.20	0.42
9:AI:83:ARG:HA	9:AI:86:VAL:HG12	2.00	0.42
10:AJ:32:ALA:N	10:AJ:78:ASN:HD21	2.17	0.42
1:AA:1123:A:C4'	10:AJ:36:GLY:HA3	2.16	0.42
12:AL:119:LYS:C	12:AL:121:GLY:N	2.73	0.42
12:AL:55:VAL:CG1	12:AL:56:ALA:N	2.80	0.42
13:AM:43:THR:HB	13:AM:44:ARG:H	1.60	0.42
14:AN:8:GLU:C	14:AN:10:ALA:N	2.72	0.42
15:AO:43:LEU:C	15:AO:45:VAL:N	2.72	0.42
16:AP:38:TYR:CE1	16:AP:50:LYS:HB3	2.55	0.42
1:AA:376:G:H5''	16:AP:5:ARG:HB2	2.01	0.42
16:AP:67:THR:CG2	16:AP:69:THR:HG23	2.49	0.42
17:AQ:6:LEU:HD23	17:AQ:6:LEU:HA	1.90	0.42
23:AW:16:C:H1'	23:AW:61:U:H1'	2.01	0.42
25:AY:131:ASN:C	25:AY:133:ARG:H	2.23	0.42
25:AY:15:GLN:HA	25:AY:15:GLN:OE1	2.20	0.42
27:B1:16:ASN:ND2	27:B1:17:SER:O	2.52	0.42
34:B8:40:GLU:CD	34:B8:44:LYS:HE3	2.39	0.42
35:BA:1034:G:C2	35:BA:1122:G:H1'	2.54	0.42
35:BA:1152:C:HO2'	51:BU:76:TYR:HE2	1.66	0.42
35:BA:1165:U:O2'	35:BA:1166:C:H5'	2.20	0.42
35:BA:1267:U:O2	35:BA:1267:U:C2'	2.65	0.42
35:BA:1300:U:C2	35:BA:1626:G:C4	3.08	0.42
33:B7:2:LYS:HG2	35:BA:1620:G:O2'	2.19	0.42
35:BA:1779:U:H5	35:BA:1784:A:N7	2.14	0.42
35:BA:1786:A:N9	35:BA:1938:A:N6	2.59	0.42
35:BA:1794:U:O2'	35:BA:1900:A:O2'	2.35	0.42
35:BA:1863:G:H1	35:BA:1879:C:H42	1.66	0.42
35:BA:1912:A:C2	35:BA:1919:A:C5	3.07	0.42
35:BA:1989:G:C2'	35:BA:1990:C:H5'	2.50	0.42
35:BA:1663:C:N3	35:BA:1992:G:O6	2.52	0.42
35:BA:2023:G:O2'	35:BA:2024:G:H5'	2.20	0.42
35:BA:218:A:O2'	35:BA:219:G:H5'	2.19	0.42
35:BA:2290:G:H1	35:BA:2342:C:N4	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2250:G:H8	35:BA:2496:C:H5''	1.84	0.42
35:BA:2508:G:C2	35:BA:2582:G:C6	3.08	0.42
35:BA:2627:G:C2	35:BA:2781:A:H2	2.37	0.42
35:BA:307:G:N2	35:BA:309:G:H3'	2.34	0.42
35:BA:438:G:H2'	35:BA:440:G:C8	2.54	0.42
35:BA:559:G:O2'	35:BA:560:C:H5'	2.19	0.42
35:BA:60:G:H21	35:BA:74:A:H2'	1.85	0.42
35:BA:826:U:H2'	35:BA:828:U:O4'	2.20	0.42
35:BA:948:G:H1'	35:BA:984:A:C2	2.54	0.42
36:BB:99:G:O2'	36:BB:100:A:O4'	2.31	0.42
38:BD:134:ARG:HB2	38:BD:135:PHE:H	1.64	0.42
38:BD:243:GLY:O	38:BD:244:ARG:HB3	2.19	0.42
38:BD:271:ILE:O	38:BD:272:ALA:O	2.36	0.42
38:BD:46:GLN:OE1	38:BD:46:GLN:N	2.52	0.42
39:BE:134:ILE:N	39:BE:134:ILE:CD1	2.80	0.42
41:BG:43:LEU:HD23	41:BG:44:GLY:N	2.24	0.42
43:BI:118:LYS:HZ2	43:BI:119:PRO:HD2	1.84	0.42
43:BI:25:TYR:CE2	43:BI:29:TYR:CD2	3.07	0.42
43:BI:28:ASN:C	43:BI:32:PRO:HG2	2.39	0.42
43:BI:80:PRO:O	43:BI:81:VAL:C	2.58	0.42
44:BN:123:TYR:CD1	44:BN:123:TYR:N	2.87	0.42
44:BN:12:ARG:NH2	44:BN:39:ARG:NH1	2.67	0.42
45:BO:43:VAL:C	45:BO:45:GLU:N	2.73	0.42
35:BA:637:A:OP2	46:BP:115:LEU:HB2	2.19	0.42
47:BQ:115:MET:CE	47:BQ:133:ARG:HH21	2.33	0.42
47:BQ:141:GLN:OE1	56:BZ:70:LEU:CB	2.68	0.42
47:BQ:54:MET:HG3	47:BQ:64:ILE:CD1	2.50	0.42
48:BR:9:LYS:O	48:BR:10:LEU:CD2	2.67	0.42
49:BS:87:PHE:HB2	49:BS:106:ARG:HD3	2.00	0.42
49:BS:28:VAL:C	49:BS:89:ARG:HG2	2.39	0.42
51:BU:21:ALA:O	51:BU:22:LYS:C	2.58	0.42
51:BU:46:ALA:O	51:BU:49:HIS:HB2	2.20	0.42
53:BW:35:ILE:HD13	53:BW:35:ILE:HA	1.88	0.42
54:BX:82:GLN:O	54:BX:85:PRO:HD2	2.19	0.42
54:BX:89:ILE:HA	54:BX:92:LEU:HB2	2.02	0.42
54:BX:88:LYS:HB3	54:BX:89:ILE:HD12	2.02	0.42
55:BY:76:CYS:CB	55:BY:77:PRO:CD	2.98	0.42
56:BZ:99:TYR:HB3	56:BZ:100:VAL:H	1.64	0.42
56:BZ:19:ARG:NH1	56:BZ:19:ARG:CB	2.82	0.42
56:BZ:6:LYS:CE	56:BZ:6:LYS:H	2.29	0.42
1:CA:1050:G:N2	1:CA:1209:C:H1'	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1054:C:OP1	1:CA:1197:G:OP2	2.37	0.42
1:CA:1225:A:H2'	1:CA:1225:A:N3	2.34	0.42
1:CA:1441:G:H5''	1:CA:1442:G:O4'	2.19	0.42
1:CA:36:C:O2'	1:CA:37:U:H5'	2.19	0.42
1:CA:544:G:C4	1:CA:545:C:C5	3.07	0.42
1:CA:751:U:H2'	1:CA:752:G:H5'	2.02	0.42
1:CA:783:C:N4	1:CA:784:C:H41	2.17	0.42
1:CA:785:G:C2'	1:CA:786:G:H5'	2.50	0.42
1:CA:868:C:H2'	1:CA:869:G:O4'	2.19	0.42
1:CA:574:A:HO2'	1:CA:882:C:HO2'	1.67	0.42
2:CB:112:VAL:HG22	2:CB:149:LEU:HD22	2.01	0.42
2:CB:168:THR:O	2:CB:169:LYS:C	2.56	0.42
1:CA:1190:G:OP2	3:CC:5:ILE:HG23	2.19	0.42
5:CE:80:ILE:CD1	5:CE:91:LEU:HB2	2.49	0.42
6:CF:100:ASN:OD1	18:CR:26:LEU:O	2.36	0.42
7:CG:15:ASP:HB3	7:CG:20:ASP:H	1.85	0.42
7:CG:16:LEU:HD11	9:CI:42:ARG:HA	2.01	0.42
11:CK:26:ASN:O	11:CK:27:ASN:CB	2.68	0.42
12:CL:76:ASN:C	12:CL:77:LEU:HD23	2.40	0.42
1:CA:551:U:O2'	12:CL:86:ARG:HD2	2.19	0.42
15:CO:82:ILE:HG23	15:CO:83:GLU:N	2.33	0.42
16:CP:73:LEU:CD2	16:CP:73:LEU:N	2.81	0.42
1:CA:277:C:OP1	17:CQ:41:LYS:HE3	2.19	0.42
18:CR:40:LEU:O	18:CR:43:PHE:N	2.52	0.42
19:CS:53:ASN:HD22	19:CS:55:LYS:H	1.66	0.42
20:CT:35:THR:O	20:CT:36:LEU:C	2.58	0.42
20:CT:93:GLU:OE1	20:CT:93:GLU:N	2.52	0.42
25:CY:72:ASP:O	25:CY:75:ALA:N	2.52	0.42
25:CY:152:ASP:OD2	26:D0:5:LYS:CB	2.67	0.42
27:D1:71:TYR:O	27:D1:74:VAL:N	2.52	0.42
34:D8:39:LYS:NZ	34:D8:40:GLU:HA	2.35	0.42
35:DA:1159:U:H2'	35:DA:1160:G:C5'	2.48	0.42
35:DA:1354:A:H2'	35:DA:1355:G:H5'	2.01	0.42
35:DA:768:G:O2'	35:DA:1379:A:N6	2.51	0.42
35:DA:1416:G:H1'	35:DA:1417:C:C6	2.55	0.42
35:DA:1619:G:C2	35:DA:1620:G:C8	3.08	0.42
35:DA:1300:U:O2	35:DA:1626:G:C4	2.72	0.42
35:DA:1711:C:H2'	35:DA:1712:C:C6	2.54	0.42
35:DA:1755:A:H2	35:DA:2716:U:C1'	2.25	0.42
35:DA:1680:U:O2	35:DA:1763:G:H3'	2.19	0.42
35:DA:1764:G:O2'	35:DA:1765:C:H5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1980:G:C2	35:DA:1982:C:C4	3.07	0.42
35:DA:2029:G:H2'	35:DA:2031:A:OP2	2.19	0.42
35:DA:2071:A:H2'	35:DA:2072:G:C8	2.54	0.42
35:DA:2186:G:H3'	35:DA:2187:G:H5''	2.01	0.42
35:DA:225:A:H2'	35:DA:226:G:O4'	2.19	0.42
35:DA:229:A:H3'	35:DA:230:U:C5'	2.43	0.42
35:DA:2407:G:C2	35:DA:2408:U:C4	3.07	0.42
35:DA:2531:A:C2	35:DA:2658:C:O2	2.73	0.42
35:DA:253:C:C2'	35:DA:254:G:H5'	2.49	0.42
35:DA:746:A:C5	35:DA:2611:U:H5''	2.54	0.42
35:DA:2736:G:C4	35:DA:2737:G:C8	3.07	0.42
35:DA:2787:C:H2'	35:DA:2787:C:O2	2.20	0.42
35:DA:2832:U:C2	35:DA:2834:G:N2	2.87	0.42
35:DA:2817:G:N2	35:DA:2836:U:H1'	2.27	0.42
35:DA:347:A:H2'	35:DA:348:G:H8	1.80	0.42
35:DA:660:G:H4'	40:DF:38:ARG:NH1	2.34	0.42
36:DB:39:A:H2'	36:DB:39:A:N3	2.35	0.42
38:DD:34:VAL:CG2	38:DD:35:LYS:HZ2	2.32	0.42
39:DE:201:THR:HG22	39:DE:203:LYS:HB3	2.01	0.42
39:DE:87:GLU:O	39:DE:88:GLY:O	2.37	0.42
40:DF:32:LEU:CD2	40:DF:33:LEU:N	2.82	0.42
41:DG:36:LYS:O	41:DG:159:VAL:HA	2.19	0.42
42:DH:19:VAL:CG1	42:DH:20:ALA:N	2.83	0.42
43:DI:111:PRO:HA	43:DI:114:LEU:CD1	2.48	0.42
43:DI:80:PRO:O	43:DI:81:VAL:C	2.58	0.42
45:DO:104:ARG:NH1	45:DO:104:ARG:CB	2.78	0.42
46:DP:80:TYR:CE1	46:DP:111:ARG:HG2	2.55	0.42
46:DP:98:GLU:HA	46:DP:98:GLU:OE1	2.20	0.42
48:DR:28:LEU:C	48:DR:30:THR:N	2.72	0.42
48:DR:59:ASP:O	48:DR:61:HIS:N	2.53	0.42
48:DR:73:VAL:HG23	48:DR:74:LYS:H	1.84	0.42
48:DR:7:GLY:O	48:DR:8:ARG:O	2.37	0.42
36:DB:7:G:C4'	49:DS:29:PHE:CE2	2.95	0.42
36:DB:7:G:C5'	49:DS:29:PHE:HE2	2.33	0.42
50:DT:108:ARG:CB	50:DT:108:ARG:HH11	2.33	0.42
50:DT:74:ARG:C	50:DT:75:ILE:HD12	2.40	0.42
51:DU:101:ARG:C	51:DU:102:GLU:HG2	2.39	0.42
53:DW:29:LEU:HD21	53:DW:33:ARG:HH21	1.84	0.42
54:DX:57:LEU:CB	54:DX:76:ARG:HD2	2.43	0.42
54:DX:81:VAL:O	54:DX:82:GLN:O	2.37	0.42
56:DZ:48:PHE:CE1	56:DZ:52:SER:HA	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:19:ARG:NH2	56:DZ:84:GLU:OE2	2.53	0.42
1:AA:1003:G:H8	1:AA:1003:G:O5'	2.03	0.42
1:AA:1063:C:H3'	1:AA:1064:G:H2'	2.00	0.42
1:AA:1188:A:H2'	1:AA:1189:C:O4'	2.19	0.42
1:AA:1192:C:P	3:AC:4:LYS:HZ1	2.43	0.42
1:AA:1293:G:HO2'	1:AA:1294:G:H8	1.61	0.42
1:AA:1420:C:H3'	1:AA:1420:C:C6	2.52	0.42
1:AA:149:A:O2'	1:AA:150:C:C6	2.69	0.42
1:AA:355:C:C2'	1:AA:356:A:H5'	2.50	0.42
1:AA:525:C:H2'	1:AA:526:C:C6	2.54	0.42
1:AA:678:U:N3	1:AA:713:G:N2	2.68	0.42
1:AA:719:C:H2'	1:AA:720:C:H5'	2.01	0.42
1:AA:780:A:C2	1:AA:803:G:C6	3.08	0.42
1:AA:826:C:H2'	1:AA:827:U:H6	1.83	0.42
1:AA:828:A:H2'	1:AA:829:G:O4'	2.19	0.42
1:AA:951:G:C2	1:AA:1231:G:C4	3.07	0.42
2:AB:142:LEU:HA	2:AB:145:LEU:HB3	2.02	0.42
2:AB:59:GLU:O	2:AB:63:MET:HG2	2.19	0.42
4:AD:10:ARG:C	4:AD:13:ARG:HB3	2.38	0.42
4:AD:63:LYS:N	4:AD:66:ARG:HH12	2.17	0.42
5:AE:73:ASN:HD22	5:AE:73:ASN:C	2.22	0.42
6:AF:21:LEU:O	6:AF:22:GLU:C	2.58	0.42
7:AG:91:VAL:CG1	7:AG:95:ARG:HD3	2.49	0.42
9:AI:45:ALA:O	9:AI:48:GLU:HB3	2.20	0.42
9:AI:66:ARG:HB3	9:AI:66:ARG:HH11	1.85	0.42
11:AK:49:GLY:C	11:AK:50:TYR:HD2	2.23	0.42
11:AK:30:VAL:HG21	11:AK:68:ALA:HB2	2.01	0.42
11:AK:79:SER:HA	11:AK:104:GLN:HB3	2.01	0.42
14:AN:34:TYR:O	14:AN:36:PHE:N	2.53	0.42
17:AQ:29:HIS:CG	17:AQ:30:PRO:HD2	2.55	0.42
19:AS:53:ASN:HD22	19:AS:55:LYS:H	1.64	0.42
25:AY:156:ARG:HH21	26:B0:6:ALA:CB	2.32	0.42
25:AY:26:ALA:C	25:AY:28:LEU:N	2.73	0.42
27:B1:13:ILE:CG1	27:B1:14:VAL:H	2.13	0.42
27:B1:37:ILE:N	27:B1:37:ILE:HD12	2.16	0.42
27:B1:38:SER:OG	27:B1:39:LYS:N	2.53	0.42
27:B1:73:LEU:HA	27:B1:76:ARG:HH11	1.80	0.42
35:BA:1005:C:H2'	35:BA:1006:C:C6	2.55	0.42
35:BA:1128:A:C2	35:BA:1129:A:C2	3.07	0.42
35:BA:1142(A):A:O2'	35:BA:1143:A:H3'	2.20	0.42
35:BA:1298:C:H2'	35:BA:1299:G:O4'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1789:A:O2'	35:BA:1790:C:H5'	2.19	0.42
35:BA:1853:A:H1'	35:BA:2233:U:O2'	2.19	0.42
35:BA:1923:U:H2'	35:BA:1924:C:H6	1.84	0.42
35:BA:2228:G:H2'	35:BA:2229:C:C6	2.55	0.42
35:BA:2287:A:N6	35:BA:2344:U:N3	2.68	0.42
35:BA:2657:A:C5	35:BA:2658:C:C5	3.08	0.42
35:BA:2679:A:C6	35:BA:2680:C:C4	3.07	0.42
35:BA:2790:A:C2'	35:BA:2790:A:N3	2.82	0.42
35:BA:2810:A:N6	35:BA:2891:G:H1'	2.35	0.42
35:BA:374:A:C2'	35:BA:375:C:H5'	2.50	0.42
35:BA:382:G:H1	35:BA:392:C:N4	2.18	0.42
35:BA:573:G:O2'	35:BA:574:C:H3'	2.19	0.42
35:BA:618:C:H2'	35:BA:619:G:H8	1.84	0.42
35:BA:61:G:H1	35:BA:94:C:N4	2.16	0.42
35:BA:954:G:OP1	47:BQ:15:GLY:N	2.48	0.42
35:BA:962:G:C2'	35:BA:963:U:H5'	2.49	0.42
36:BB:46:A:H2'	36:BB:47:C:C6	2.55	0.42
35:BA:2123:G:H1'	37:BC:175:VAL:CB	2.50	0.42
38:BD:96:HIS:ND1	38:BD:102:LYS:HD2	2.35	0.42
40:BF:132:VAL:CG2	40:BF:133:ASN:N	2.71	0.42
40:BF:139:PHE:HB2	40:BF:166:ALA:HB1	2.00	0.42
41:BG:107:LEU:HD23	41:BG:111:LEU:CD1	2.49	0.42
41:BG:174:GLU:C	41:BG:176:LEU:N	2.73	0.42
41:BG:4:ASP:O	41:BG:5:VAL:HG22	2.20	0.42
42:BH:105:LEU:HD22	42:BH:105:LEU:H	1.84	0.42
42:BH:122:THR:CB	42:BH:134:SER:HB2	2.44	0.42
43:BI:94:ALA:HB1	43:BI:114:LEU:HD12	2.00	0.42
43:BI:132:PRO:O	43:BI:135:GLU:HG2	2.20	0.42
44:BN:22:THR:N	44:BN:61:ARG:HB2	2.34	0.42
44:BN:96:GLU:N	44:BN:96:GLU:OE2	2.50	0.42
46:BP:19:VAL:C	46:BP:20:GLY:O	2.58	0.42
35:BA:1242:A:N1	46:BP:8:PRO:CG	2.82	0.42
47:BQ:54:MET:CG	47:BQ:64:ILE:HD13	2.50	0.42
48:BR:28:LEU:HD12	48:BR:29:LEU:HD13	2.01	0.42
50:BT:101:PHE:C	50:BT:101:PHE:HD2	2.23	0.42
35:BA:2864:G:OP1	50:BT:119:LYS:HD2	2.19	0.42
50:BT:52:ILE:HA	50:BT:61:PHE:HA	2.01	0.42
51:BU:111:GLU:O	51:BU:112:ARG:C	2.57	0.42
51:BU:57:PHE:CD1	51:BU:60:LEU:HD12	2.54	0.42
52:BV:78:LYS:C	52:BV:78:LYS:CD	2.88	0.42
53:BW:107:LEU:HA	53:BW:107:LEU:HD12	1.79	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:BZ:129:SER:HA	56:BZ:130:PRO:HD3	1.87	0.42
56:BZ:85:HIS:ND1	56:BZ:86:VAL:N	2.68	0.42
1:CA:227:G:C6	1:CA:228:A:C6	3.08	0.42
1:CA:246:A:O3'	1:CA:247:G:H4'	2.19	0.42
1:CA:287:U:O2'	1:CA:288:A:H5'	2.19	0.42
1:CA:435:C:N3	1:CA:436:C:C5	2.88	0.42
1:CA:525:C:H2'	1:CA:526:C:C6	2.55	0.42
1:CA:611:A:C2'	1:CA:612:C:H5'	2.50	0.42
1:CA:66:G:H4'	1:CA:173:U:C4	2.52	0.42
1:CA:699:C:H2'	1:CA:700:G:H8	1.84	0.42
2:CB:153:ARG:O	2:CB:154:LEU:O	2.37	0.42
3:CC:206:GLU:O	3:CC:207:VAL:C	2.57	0.42
3:CC:22:TRP:HZ2	3:CC:36:ASP:OD1	2.02	0.42
3:CC:58:GLU:HB2	3:CC:65:ALA:HB2	2.01	0.42
1:CA:688:G:H5'	11:CK:46:GLY:C	2.39	0.42
11:CK:60:ALA:C	11:CK:62:GLN:N	2.72	0.42
13:CM:48:LEU:HD21	13:CM:53:VAL:CG2	2.49	0.42
13:CM:83:ASP:CG	13:CM:84:ILE:H	2.23	0.42
15:CO:11:VAL:O	15:CO:14:GLU:HB3	2.19	0.42
16:CP:40:ASP:OD2	16:CP:40:ASP:C	2.56	0.42
16:CP:45:THR:O	16:CP:46:PRO:C	2.58	0.42
16:CP:55:ARG:O	16:CP:58:TYR:N	2.52	0.42
23:CW:67:C:H2'	23:CW:68:C:C6	2.55	0.42
25:CY:175:LEU:O	25:CY:178:LYS:HB2	2.20	0.42
25:CY:53:ASN:OD1	25:CY:54:GLN:N	2.43	0.42
26:D0:51:VAL:HG13	26:D0:60:PHE:O	2.19	0.42
26:D0:55:ARG:C	26:D0:57:PHE:H	2.23	0.42
26:D0:56:ASP:O	26:D0:58:THR:N	2.53	0.42
29:D3:51:ALA:C	29:D3:53:LEU:H	2.23	0.42
33:D7:8:ASN:HD22	33:D7:9:ARG:H	1.59	0.42
35:DA:1210:A:H1'	35:DA:1212:G:N3	2.33	0.42
35:DA:1257:C:H2'	35:DA:1258:C:C6	2.55	0.42
35:DA:1323:U:H3	35:DA:1331:A:N6	2.13	0.42
35:DA:1938:A:H4'	35:DA:1939:U:OP2	2.20	0.42
35:DA:563:G:C6	35:DA:2018:G:C5	3.08	0.42
35:DA:214:G:O2'	35:DA:215:G:P	2.77	0.42
35:DA:2291:U:OP1	35:DA:2381:C:H5'	2.19	0.42
35:DA:2059:A:C5	35:DA:2503:A:C2	3.07	0.42
35:DA:2866:U:C6	35:DA:2868:A:C1'	3.02	0.42
35:DA:708:C:H5'	35:DA:709:U:OP2	2.20	0.42
35:DA:58:G:N3	35:DA:70:G:C2	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:D5:3:LYS:HB2	35:DA:747:U:C5	2.54	0.42
26:D0:77:ARG:NH2	35:DA:857:C:OP2	2.53	0.42
35:DA:936:C:H2'	35:DA:937:U:C6	2.54	0.42
38:DD:166:GLN:HB2	38:DD:174:ILE:HG22	2.01	0.42
38:DD:265:PRO:O	38:DD:266:SER:C	2.58	0.42
39:DE:33:VAL:CG1	39:DE:89:ASP:N	2.73	0.42
40:DF:185:ASP:OD1	40:DF:188:ARG:HD2	2.20	0.42
41:DG:107:LEU:HD23	41:DG:111:LEU:CD1	2.49	0.42
41:DG:165:THR:OG1	41:DG:168:GLU:OE1	2.30	0.42
41:DG:28:VAL:HG12	41:DG:28:VAL:O	2.19	0.42
43:DI:102:SER:HB2	43:DI:109:ILE:HG23	2.01	0.42
44:DN:79:PRO:HG2	44:DN:80:GLY:N	2.34	0.42
44:DN:99:LEU:HD22	44:DN:99:LEU:HA	1.84	0.42
45:DO:62:VAL:O	45:DO:63:VAL:CG1	2.67	0.42
35:DA:389:G:N1	46:DP:71:VAL:CG1	2.81	0.42
47:DQ:55:VAL:C	47:DQ:57:HIS:N	2.72	0.42
50:DT:119:LYS:O	50:DT:123:GLN:HG2	2.19	0.42
45:DO:104:ARG:NH2	50:DT:33:LYS:CE	2.82	0.42
50:DT:50:ILE:HD11	50:DT:64:ARG:HB3	2.02	0.42
51:DU:103:PRO:C	51:DU:105:VAL:N	2.71	0.42
51:DU:111:GLU:O	51:DU:115:ALA:CB	2.67	0.42
51:DU:7:GLY:O	51:DU:8:VAL:CG2	2.68	0.42
51:DU:83:LEU:HD12	51:DU:113:ALA:HB2	2.00	0.42
51:DU:95:LEU:HD23	51:DU:95:LEU:HA	1.88	0.42
54:DX:23:GLU:HG3	54:DX:24:GLY:N	2.34	0.42
54:DX:61:GLY:H	54:DX:70:LEU:CD2	2.32	0.42
54:DX:88:LYS:HB3	54:DX:89:ILE:CD1	2.49	0.42
55:DY:71:LYS:HZ2	55:DY:71:LYS:HB2	1.84	0.42
56:DZ:58:VAL:O	56:DZ:59:LEU:C	2.58	0.42
1:AA:106:C:O2	1:AA:379:C:H4'	2.20	0.42
1:AA:107:G:C2'	1:AA:108:G:H5'	2.50	0.42
1:AA:119:A:OP2	1:AA:288:A:N6	2.50	0.42
1:AA:1235:U:O3'	21:AU:3:LYS:HB2	2.20	0.42
1:AA:1473:A:H2'	1:AA:1474:G:C8	2.55	0.42
1:AA:1402:C:O2	1:AA:1500:A:N1	2.52	0.42
1:AA:287:U:O2'	1:AA:288:A:H5'	2.19	0.42
1:AA:377:G:OP1	16:AP:5:ARG:NH1	2.50	0.42
1:AA:446:G:C2'	1:AA:447:G:H5'	2.49	0.42
1:AA:545:C:OP1	4:AD:61:LYS:NZ	2.53	0.42
1:AA:300:A:H2	1:AA:566:G:O6	2.03	0.42
1:AA:630:G:H2'	1:AA:631:G:H5'	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:778:G:H2'	1:AA:779:C:O4'	2.20	0.42
1:AA:961:U:C2'	1:AA:962:C:H5'	2.50	0.42
1:AA:965:A:OP1	1:AA:1198:G:H5''	2.19	0.42
4:AD:15:GLU:N	4:AD:15:GLU:CD	2.73	0.42
4:AD:162:LEU:O	4:AD:163:GLU:O	2.38	0.42
4:AD:176:LEU:HD12	4:AD:177:ASP:N	2.33	0.42
5:AE:78:HIS:HB2	5:AE:79:GLU:H	1.69	0.42
8:AH:120:THR:HG23	8:AH:123:GLU:CD	2.40	0.42
8:AH:12:ARG:HA	8:AH:15:ASN:HD22	1.84	0.42
9:AI:4:TYR:H	9:AI:4:TYR:HD1	1.68	0.42
11:AK:59:TYR:CE2	11:AK:63:LEU:HD11	2.53	0.42
11:AK:64:ALA:C	11:AK:66:LEU:N	2.73	0.42
12:AL:48:PRO:HD2	12:AL:49:ASN:OD1	2.19	0.42
17:AQ:27:PHE:C	17:AQ:27:PHE:CD1	2.92	0.42
19:AS:78:ARG:N	19:AS:78:ARG:CD	2.83	0.42
20:AT:38:LYS:HE2	20:AT:38:LYS:HB3	1.84	0.42
22:AV:34:A:O2'	22:AV:35:A:H5'	2.20	0.42
23:AW:72:C:H6	23:AW:72:C:O5'	2.02	0.42
25:AY:175:LEU:C	25:AY:177:GLU:N	2.71	0.42
25:AY:28:LEU:HB3	25:AY:114:LEU:CD1	2.49	0.42
29:B3:16:PRO:CB	29:B3:18:ASP:OD1	2.67	0.42
29:B3:17:LYS:O	29:B3:18:ASP:C	2.58	0.42
29:B3:1:MET:O	29:B3:3:ARG:N	2.51	0.42
34:B8:39:LYS:NZ	34:B8:40:GLU:HA	2.35	0.42
35:BA:1251:C:OP2	35:BA:1251:C:H2'	2.20	0.42
35:BA:1277:G:H2'	35:BA:1278:A:O4'	2.19	0.42
35:BA:1406:U:C3'	35:BA:1407:C:H6	2.30	0.42
35:BA:1514:U:C2	35:BA:1515:G:C8	3.08	0.42
35:BA:1714:G:H2'	35:BA:1717:G:C8	2.53	0.42
35:BA:961:C:C5	35:BA:2031:A:C2	3.08	0.42
35:BA:2040:C:H2'	35:BA:2041:U:H6	1.82	0.42
35:BA:214:G:O2'	35:BA:215:G:P	2.77	0.42
35:BA:2264:C:H2'	35:BA:2265:U:O4'	2.19	0.42
35:BA:2415:G:H2'	35:BA:2416:C:C6	2.55	0.42
35:BA:2491:U:O2'	35:BA:2492:U:H5'	2.20	0.42
35:BA:2737:G:N3	35:BA:2738:A:C8	2.88	0.42
35:BA:363(F):A:O2'	35:BA:364:C:C5	2.71	0.42
35:BA:571:A:C5'	35:BA:2030:A:N6	2.73	0.42
35:BA:702:G:N3	35:BA:731:C:N3	2.67	0.42
28:B2:55:ARG:CZ	35:BA:72:U:OP1	2.67	0.42
35:BA:790:C:OP1	35:BA:790:C:H4'	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:949:C:O2'	35:BA:950:G:H5'	2.20	0.42
36:BB:6:C:C4	36:BB:116:G:N1	2.88	0.42
38:BD:147:LEU:HB3	38:BD:148:GLU:H	1.59	0.42
38:BD:177:LEU:HB3	38:BD:178:PRO:CD	2.50	0.42
38:BD:109:ASP:HB2	38:BD:197:GLY:HA2	2.01	0.42
38:BD:257:LEU:HD23	38:BD:258:LYS:O	2.19	0.42
40:BF:41:LEU:HD11	40:BF:184:TYR:CE1	2.48	0.42
41:BG:60:LEU:O	41:BG:63:ILE:CG1	2.55	0.42
43:BI:77:LEU:CD2	43:BI:101:LEU:HD13	2.50	0.42
44:BN:115:ARG:O	44:BN:118:LYS:HB2	2.20	0.42
34:B8:25:MET:CG	46:BP:64:LYS:HB2	2.50	0.42
47:BQ:55:VAL:HG22	47:BQ:56:ARG:N	2.35	0.42
47:BQ:77:LYS:HA	47:BQ:78:PRO:HD3	1.84	0.42
48:BR:103:ARG:NH1	48:BR:103:ARG:CG	2.81	0.42
48:BR:20:LEU:CD1	48:BR:21:TYR:N	2.83	0.42
49:BS:28:VAL:H	49:BS:89:ARG:CG	2.33	0.42
51:BU:103:PRO:C	51:BU:105:VAL:N	2.73	0.42
51:BU:52:ARG:O	51:BU:55:ARG:N	2.52	0.42
53:BW:19:LEU:HA	53:BW:19:LEU:HD12	1.84	0.42
53:BW:29:LEU:HD21	53:BW:33:ARG:NH2	2.34	0.42
56:BZ:3:TYR:O	56:BZ:57:ILE:CG2	2.68	0.42
56:BZ:97:GLU:O	56:BZ:98:MET:CB	2.66	0.42
1:CA:1009:G:O2'	1:CA:1010:G:H5'	2.20	0.42
1:CA:106:C:H2'	1:CA:107:G:H8	1.85	0.42
1:CA:1346:A:O3'	1:CA:1347:G:H4'	2.19	0.42
1:CA:1358:U:H2'	1:CA:1359:C:O4'	2.20	0.42
1:CA:1371:G:OP2	9:CI:11:LYS:HE3	2.19	0.42
1:CA:1387:G:H2'	1:CA:1388:C:C6	2.55	0.42
1:CA:1486:G:C6	1:CA:1487:G:N1	2.88	0.42
1:CA:267:C:OP2	17:CQ:67:LYS:HD2	2.19	0.42
1:CA:594:G:H2'	1:CA:595:G:O4'	2.19	0.42
1:CA:73:G:C2	1:CA:97:G:N1	2.88	0.42
1:CA:766:A:H62	1:CA:813:U:H3	1.66	0.42
1:CA:901:A:C5	1:CA:902:G:H1'	2.54	0.42
2:CB:116:GLU:HG2	2:CB:116:GLU:H	1.56	0.42
3:CC:23:TYR:HA	10:CJ:11:PHE:HE1	1.81	0.42
4:CD:172:PRO:C	4:CD:187:ARG:HH12	2.23	0.42
5:CE:36:ASP:O	5:CE:37:ARG:CB	2.65	0.42
7:CG:64:GLN:HG2	7:CG:128:ALA:HB1	2.01	0.42
7:CG:15:ASP:OD2	7:CG:44:TYR:OH	2.36	0.42
1:CA:1240:U:N3	7:CG:30:ILE:HG22	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CJ:8:LEU:CA	10:CJ:96:ILE:HG22	2.49	0.42
11:CK:21:ILE:HG13	11:CK:30:VAL:HG12	2.01	0.42
13:CM:120:LYS:CE	13:CM:120:LYS:HA	2.41	0.42
16:CP:60:LEU:H	16:CP:60:LEU:HG	1.61	0.42
19:CS:42:PRO:O	19:CS:44:MET:N	2.47	0.42
25:CY:10:THR:HG22	25:CY:14:MET:SD	2.59	0.42
25:CY:159:ALA:O	25:CY:162:GLN:HB3	2.18	0.42
25:CY:32:ARG:HA	25:CY:32:ARG:CZ	2.50	0.42
27:D1:38:SER:O	27:D1:39:LYS:HB3	2.18	0.42
28:D2:14:ARG:NE	28:D2:15:LYS:N	2.68	0.42
35:DA:1158:C:O2'	35:DA:1159:U:H6	2.02	0.42
35:DA:1266:G:H22	35:DA:2012:G:H2'	1.84	0.42
35:DA:1279:G:H2'	35:DA:1280:G:C8	2.54	0.42
35:DA:1281:G:O6	35:DA:1286:A:N7	2.52	0.42
35:DA:1417:C:H2'	35:DA:1418:G:H5'	2.01	0.42
35:DA:1426:G:C6	35:DA:1427:A:N1	2.87	0.42
35:DA:1444:G:H2'	35:DA:1445(A):C:C5	2.55	0.42
35:DA:1629:U:H2'	35:DA:1630:G:C8	2.55	0.42
35:DA:1635:G:O2'	35:DA:1636:C:H5'	2.19	0.42
35:DA:1658:C:H2'	35:DA:1659:U:H6	1.84	0.42
35:DA:1796:U:H2'	35:DA:1797:C:C6	2.55	0.42
35:DA:1819:A:H1'	35:DA:1821:A:C5	2.55	0.42
35:DA:2026:C:H2'	35:DA:2027:G:O4'	2.20	0.42
35:DA:2114:A:C3'	35:DA:2115:G:H5'	2.49	0.42
35:DA:214:G:C2'	35:DA:215:G:OP2	2.67	0.42
35:DA:2173:A:H3'	35:DA:2173:A:P	2.60	0.42
35:DA:2183:C:O2'	35:DA:2184:G:H5'	2.19	0.42
35:DA:2292:C:H2'	35:DA:2293:C:H6	1.84	0.42
35:DA:2415:G:H4'	46:DP:67:MET:N	2.33	0.42
35:DA:511:U:O4	35:DA:512:G:C2	2.72	0.42
35:DA:666:G:O2'	35:DA:667:U:H5'	2.19	0.42
35:DA:740:U:C2	35:DA:758:C:H1'	2.55	0.42
35:DA:856:C:H6	35:DA:856:C:H5''	1.84	0.42
38:DD:267:SER:HA	38:DD:270:ILE:CG1	2.49	0.42
39:DE:1:MET:N	39:DE:84:PHE:HB2	2.34	0.42
40:DF:116:ASP:OD2	46:DP:5:ASP:N	2.53	0.42
40:DF:160:ASN:ND2	40:DF:160:ASN:C	2.72	0.42
40:DF:3:GLU:HB3	40:DF:20:LEU:O	2.19	0.42
40:DF:65:TRP:HZ3	40:DF:73:ALA:O	2.03	0.42
41:DG:125:PHE:CE2	41:DG:131:TYR:HD2	2.37	0.42
44:DN:46:VAL:HG22	44:DN:47:ALA:H	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:55:VAL:CG1	44:DN:56:ASN:N	2.82	0.42
45:DO:13:ASN:ND2	45:DO:96:THR:OG1	2.53	0.42
45:DO:37:ASP:HB2	45:DO:62:VAL:HG23	2.01	0.42
45:DO:98:VAL:CG2	45:DO:99:PHE:N	2.82	0.42
47:DQ:23:GLY:CA	47:DQ:101:ARG:HB2	2.46	0.42
47:DQ:6:ARG:O	47:DQ:6:ARG:HG3	2.20	0.42
47:DQ:12:GLN:HE21	47:DQ:72:LYS:HA	1.85	0.42
47:DQ:93:TYR:N	47:DQ:93:TYR:CD1	2.86	0.42
48:DR:2:ARG:CZ	48:DR:5:LYS:CE	2.97	0.42
50:DT:53:ARG:HG2	50:DT:53:ARG:NH1	2.26	0.42
51:DU:54:LYS:C	51:DU:56:ASP:H	2.23	0.42
51:DU:70:ARG:O	51:DU:73:GLY:N	2.43	0.42
52:DV:57:VAL:O	52:DV:57:VAL:HG12	2.20	0.42
55:DY:31:LEU:CG	55:DY:34:LYS:HB2	2.50	0.42
55:DY:42:VAL:CB	55:DY:65:ALA:HB3	2.49	0.42
56:DZ:137:ILE:CD1	56:DZ:157:LEU:HA	2.49	0.42
56:DZ:48:PHE:O	56:DZ:49:ARG:C	2.58	0.42
1:AA:1071:C:H2'	1:AA:1072:G:H8	1.84	0.42
1:AA:1132:C:C2'	1:AA:1133:G:H5'	2.50	0.42
1:AA:1191:A:H5''	3:AC:4:LYS:HZ2	1.84	0.42
1:AA:1320:C:H5'	19:AS:70:LYS:HD3	2.02	0.42
1:AA:434:U:H2'	1:AA:435:C:N1	2.33	0.42
1:AA:686:U:C5	1:AA:687:A:N6	2.87	0.42
1:AA:69:G:H2'	1:AA:70:G:C8	2.55	0.42
1:AA:783:C:C6	1:AA:784:C:H5	2.36	0.42
1:AA:929:G:O2'	1:AA:930:C:H5'	2.19	0.42
1:AA:1074:G:O2'	2:AB:103:THR:HG21	2.20	0.42
4:AD:58:LEU:C	4:AD:58:LEU:HD13	2.40	0.42
5:AE:126:ARG:HG3	5:AE:126:ARG:HH11	1.85	0.42
5:AE:60:TYR:CE2	5:AE:64:ARG:NH2	2.84	0.42
5:AE:76:ILE:CG2	5:AE:77:PRO:N	2.83	0.42
9:AI:54:ASP:O	9:AI:58:ARG:HB2	2.20	0.42
10:AJ:78:ASN:O	10:AJ:80:LYS:N	2.53	0.42
11:AK:107:SER:C	11:AK:108:ILE:HD12	2.40	0.42
17:AQ:63:ARG:HG2	17:AQ:64:PRO:HD2	2.02	0.42
18:AR:53:ARG:O	18:AR:54:ARG:C	2.58	0.42
18:AR:73:ALA:HB1	18:AR:78:LEU:HB2	2.01	0.42
1:AA:261:U:OP2	20:AT:79:ARG:NH2	2.53	0.42
25:AY:150:SER:OG	25:AY:153:GLU:CG	2.65	0.42
35:BA:1208:C:C5	35:BA:1209:G:N7	2.88	0.42
35:BA:1321:A:H2'	35:BA:1321:A:N3	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1381:G:O2'	35:BA:1382:G:H5'	2.19	0.42
35:BA:139(A):G:H5'	35:BA:140:G:OP2	2.19	0.42
35:BA:1544:A:H2'	35:BA:1545:A:OP2	2.19	0.42
35:BA:1615:C:C5	35:BA:1617:C:C2	3.08	0.42
35:BA:171:G:C5	35:BA:172:C:C5	3.07	0.42
35:BA:1681:G:H1'	35:BA:1763:G:H5'	2.01	0.42
35:BA:1826:G:H4'	38:BD:242:ARG:CZ	2.49	0.42
35:BA:2128:C:H5'	35:BA:2173:A:H2	1.85	0.42
35:BA:2206:G:N1	35:BA:2208:A:OP1	2.53	0.42
35:BA:2725:A:N7	35:BA:2727:G:C5	2.87	0.42
35:BA:2809:A:C2'	35:BA:2810:A:H5'	2.50	0.42
35:BA:2863:C:C3'	35:BA:2864:G:C5'	2.98	0.42
35:BA:395:U:H1'	35:BA:396:G:N7	2.35	0.42
35:BA:98:G:N3	35:BA:98:G:H2'	2.34	0.42
36:BB:40:U:C2'	36:BB:41:U:OP1	2.67	0.42
37:BC:59:ARG:HB2	37:BC:62:VAL:CG2	2.39	0.42
37:BC:59:ARG:O	37:BC:62:VAL:HG22	2.20	0.42
38:BD:203:ASN:O	38:BD:204:ILE:C	2.57	0.42
38:BD:270:ILE:O	38:BD:271:ILE:HG13	2.20	0.42
38:BD:43:ARG:HB3	38:BD:54:ARG:HB2	2.01	0.42
41:BG:46:ALA:CA	41:BG:51:ARG:HG3	2.50	0.42
41:BG:7:LEU:O	41:BG:11:TYR:N	2.46	0.42
43:BI:115:ALA:HB2	43:BI:129:THR:O	2.19	0.42
44:BN:18:ALA:C	44:BN:20:GLY:N	2.73	0.42
44:BN:44:PRO:C	44:BN:46:VAL:N	2.72	0.42
45:BO:115:VAL:O	45:BO:118:ALA:HB3	2.19	0.42
45:BO:11:ALA:O	45:BO:98:VAL:HG23	2.20	0.42
46:BP:18:ARG:HE	46:BP:18:ARG:HB3	1.58	0.42
46:BP:7:ARG:HA	46:BP:7:ARG:HD2	1.85	0.42
48:BR:31:HIS:O	48:BR:33:ARG:N	2.52	0.42
48:BR:54:LEU:HD22	48:BR:66:VAL:HG22	2.01	0.42
49:BS:54:LEU:O	49:BS:57:LYS:N	2.52	0.42
49:BS:77:ALA:O	49:BS:79:ALA:N	2.52	0.42
50:BT:58:ASN:ND2	50:BT:58:ASN:H	2.17	0.42
50:BT:83:ILE:O	50:BT:84:GLN:C	2.58	0.42
53:BW:20:VAL:O	53:BW:21:VAL:C	2.58	0.42
53:BW:46:PHE:C	53:BW:48:ALA:N	2.71	0.42
53:BW:65:LEU:HD21	53:BW:67:ASP:HB2	2.01	0.42
53:BW:86:LEU:HA	53:BW:87:PRO:HD3	1.83	0.42
54:BX:15:GLU:O	54:BX:19:ALA:N	2.45	0.42
54:BX:82:GLN:CD	54:BX:83:VAL:HG22	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:28:LYS:O	55:BY:37:VAL:C	2.58	0.42
55:BY:8:LYS:HD2	55:BY:8:LYS:N	2.20	0.42
56:BZ:99:TYR:O	56:BZ:100:VAL:CB	2.68	0.42
1:CA:1060:C:H4'	10:CJ:52:GLY:CA	2.50	0.42
1:CA:1148:U:H2'	1:CA:1149:C:O4'	2.20	0.42
1:CA:1117:G:N2	1:CA:1180:A:H1'	2.35	0.42
1:CA:1277:C:H4'	1:CA:1282:C:O2	2.20	0.42
1:CA:1443:G:C6	1:CA:1444:C:N4	2.88	0.42
1:CA:1508:G:C4	1:CA:1509:C:C6	3.07	0.42
1:CA:1518:A:C2	1:CA:1519:A:N1	2.88	0.42
1:CA:176:C:C2	1:CA:177:C:C5	3.08	0.42
1:CA:29:G:H5'	1:CA:296:U:OP1	2.19	0.42
1:CA:60:A:H4'	1:CA:61:G:O5'	2.19	0.42
1:CA:640:A:H2'	1:CA:641:U:H5'	2.00	0.42
1:CA:68:G:N2	1:CA:69:G:H1'	2.35	0.42
1:CA:709:G:C4	1:CA:710:G:C8	3.08	0.42
1:CA:745:C:O2'	1:CA:746:A:H5'	2.19	0.42
1:CA:830:G:C6	1:CA:831:U:N3	2.87	0.42
2:CB:101:MET:HB2	2:CB:102:LEU:CD1	2.47	0.42
2:CB:17:PHE:O	2:CB:18:GLY:O	2.38	0.42
1:CA:1103:C:H5''	2:CB:98:LEU:CD1	2.50	0.42
4:CD:105:VAL:HG13	4:CD:110:PHE:HB2	2.02	0.42
4:CD:45:GLN:O	4:CD:46:LYS:HG3	2.20	0.42
6:CF:67:MET:HB2	6:CF:68:PRO:CD	2.48	0.42
8:CH:60:ARG:HG3	8:CH:60:ARG:NH1	2.34	0.42
1:CA:1228:C:OP1	13:CM:115:LYS:HG3	2.19	0.42
13:CM:36:LYS:HB2	13:CM:59:TYR:CZ	2.55	0.42
13:CM:91:ARG:CB	13:CM:96:LEU:O	2.65	0.42
14:CN:29:ARG:HH12	14:CN:31:ARG:CG	2.33	0.42
3:CC:37:GLN:NE2	14:CN:52:GLN:OE1	2.53	0.42
15:CO:64:ARG:O	15:CO:66:LEU:N	2.53	0.42
16:CP:2:VAL:O	16:CP:2:VAL:HG22	2.20	0.42
16:CP:34:GLU:OE1	16:CP:36:ILE:HG23	2.20	0.42
16:CP:67:THR:CG2	16:CP:69:THR:HG23	2.49	0.42
18:CR:76:LEU:C	18:CR:78:LEU:H	2.23	0.42
1:CA:1320:C:P	19:CS:70:LYS:HE3	2.60	0.42
25:CY:151:GLU:O	25:CY:151:GLU:HG2	2.20	0.42
27:D1:20:ARG:HG2	27:D1:20:ARG:H	1.46	0.42
29:D3:12:PRO:HA	29:D3:15:TYR:HD1	1.84	0.42
31:D5:13:LYS:HZ1	35:DA:516:C:P	2.43	0.42
32:D6:47:THR:CG2	32:D6:48:VAL:N	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:D8:22:VAL:HB	34:D8:53:PRO:HB3	1.97	0.42
35:DA:1004:C:O4'	35:DA:1010:A:C6	2.73	0.42
35:DA:1213:A:H2'	35:DA:1214:A:C8	2.48	0.42
35:DA:1560:G:H2'	35:DA:1561:G:H8	1.85	0.42
35:DA:1313:U:H2'	35:DA:1610:A:C2	2.54	0.42
35:DA:1793:C:O2'	35:DA:1794:U:H5'	2.20	0.42
35:DA:2206:G:N1	35:DA:2208:A:OP1	2.53	0.42
35:DA:2220:G:C4	35:DA:2221:G:C8	3.07	0.42
35:DA:966:G:C1'	35:DA:2267:A:H62	2.33	0.42
35:DA:2059:A:H62	35:DA:2503:A:H2'	1.85	0.42
35:DA:2513:G:C6	35:DA:2514:U:C4	3.08	0.42
35:DA:271(L):U:H5"	35:DA:271(M):G:C4	2.54	0.42
35:DA:2744:G:N2	35:DA:2761:G:C4	2.88	0.42
35:DA:339:U:O5'	35:DA:339:U:H6	2.03	0.42
35:DA:656:G:C6	35:DA:657:U:C4	3.08	0.42
35:DA:664:C:H4'	35:DA:941:A:OP1	2.20	0.42
35:DA:671:C:C5	46:DP:36:LYS:NZ	2.86	0.42
35:DA:785:G:C6	35:DA:786:C:C4	3.08	0.42
35:DA:870:A:C2	35:DA:871:U:H1'	2.55	0.42
36:DB:7:G:H5'	36:DB:8:U:OP2	2.20	0.42
37:DC:18:LYS:O	37:DC:20:TYR:N	2.52	0.42
38:DD:257:LEU:HD23	38:DD:258:LYS:C	2.40	0.42
39:DE:71:GLY:O	39:DE:72:VAL:HB	2.20	0.42
40:DF:25:PRO:HB3	40:DF:119:ARG:CD	2.50	0.42
41:DG:133:LEU:N	41:DG:133:LEU:HD12	2.34	0.42
41:DG:142:PRO:HG2	41:DG:143:GLU:N	2.34	0.42
41:DG:146:TYR:HD2	41:DG:146:TYR:O	2.03	0.42
43:DI:118:LYS:HB3	43:DI:118:LYS:HE3	1.88	0.42
35:DA:6:A:C2'	44:DN:130:HIS:HB2	2.49	0.42
45:DO:24:VAL:O	45:DO:24:VAL:HG13	2.20	0.42
46:DP:56:SER:O	46:DP:57:THR:CB	2.66	0.42
47:DQ:134:ARG:CG	47:DQ:135:ASP:N	2.82	0.42
48:DR:53:HIS:O	48:DR:56:LYS:HB2	2.19	0.42
49:DS:43:GLU:HB2	49:DS:44:LYS:H	1.63	0.42
49:DS:73:LEU:O	49:DS:74:ALA:C	2.58	0.42
35:DA:2864:G:OP1	50:DT:119:LYS:HD2	2.20	0.42
45:DO:78:ARG:HB3	50:DT:73:GLU:CG	2.50	0.42
52:DV:78:LYS:CD	52:DV:78:LYS:C	2.88	0.42
53:DW:65:LEU:HD21	53:DW:67:ASP:HB2	2.02	0.42
54:DX:62:LYS:CB	54:DX:68:ARG:HB2	2.41	0.42
54:DX:78:LYS:HD3	54:DX:78:LYS:C	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:62:GLU:OE1	55:DY:62:GLU:CA	2.68	0.42
55:DY:86:ARG:CB	55:DY:88:LYS:HZ2	2.25	0.42
56:DZ:128:VAL:HG22	56:DZ:129:SER:N	2.34	0.42
1:AA:1204:A:N1	1:AA:1205:U:O2	2.53	0.42
1:AA:1206:G:H1'	3:AC:193:TYR:O	2.20	0.42
1:AA:1299:A:C8	1:AA:1301:U:C2	3.08	0.42
1:AA:1350:A:C2	1:AA:1351:U:C2	3.07	0.42
1:AA:1440:C:H2'	1:AA:1441:G:C8	2.55	0.42
1:AA:1471:G:O2'	1:AA:1472:U:H5'	2.19	0.42
1:AA:259:G:H2'	1:AA:260:G:H8	1.83	0.42
1:AA:381:C:H2'	1:AA:382:A:O4'	2.20	0.42
1:AA:502:G:C6	1:AA:544:G:C6	3.07	0.42
1:AA:652:U:C2	1:AA:752:G:N2	2.88	0.42
1:AA:709:G:C4	1:AA:710:G:C8	3.07	0.42
1:AA:745:C:O2'	1:AA:746:A:H5'	2.20	0.42
1:AA:769:G:H1	1:AA:810:C:N4	2.18	0.42
1:AA:783:C:N4	1:AA:800:G:N2	2.67	0.42
1:AA:821:G:O2'	1:AA:822:C:H5'	2.20	0.42
1:AA:835:U:H3	1:AA:851:G:H1	1.67	0.42
2:AB:114:ARG:HH11	2:AB:118:LEU:CD2	2.33	0.42
1:AA:1100:C:OP2	2:AB:96:ARG:HG2	2.20	0.42
1:AA:1103:C:H5''	2:AB:98:LEU:CD1	2.50	0.42
3:AC:150:LYS:HA	3:AC:168:ALA:O	2.19	0.42
3:AC:22:TRP:HZ2	3:AC:36:ASP:OD1	2.03	0.42
3:AC:60:ALA:O	3:AC:61:ALA:HB2	2.20	0.42
4:AD:62:GLN:HB3	4:AD:66:ARG:HH12	1.85	0.42
7:AG:122:HIS:O	7:AG:123:GLU:C	2.58	0.42
8:AH:26:VAL:CG2	8:AH:32:LYS:HZ3	2.29	0.42
10:AJ:23:ILE:HG23	10:AJ:85:LEU:HD22	2.02	0.42
1:AA:657:G:N2	15:AO:22:THR:OG1	2.53	0.42
15:AO:55:GLY:O	15:AO:59:MET:HG3	2.19	0.42
17:AQ:67:LYS:C	17:AQ:69:LYS:H	2.23	0.42
20:AT:57:ARG:HB3	20:AT:58:LYS:H	1.71	0.42
20:AT:83:ARG:HG2	20:AT:86:ARG:HD3	2.01	0.42
23:AW:16:C:O4'	23:AW:60:A:C2	2.72	0.42
23:AW:57:C:H2'	23:AW:58:A:C8	2.51	0.42
25:AY:127:VAL:O	25:AY:128:ALA:C	2.58	0.42
25:AY:156:ARG:NH2	47:BQ:80:GLU:CB	2.82	0.42
25:AY:64:ARG:HA	25:AY:103:ILE:CG1	2.50	0.42
25:AY:67:VAL:O	25:AY:67:VAL:HG23	2.20	0.42
26:B0:55:ARG:C	26:B0:57:PHE:H	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:B5:15:ARG:HG3	31:B5:15:ARG:NH1	2.34	0.42
34:B8:4:MET:HE2	34:B8:61:LEU:HD12	2.01	0.42
34:B8:22:VAL:CB	34:B8:53:PRO:HB2	2.37	0.42
34:B8:11:LYS:HA	34:B8:60:LEU:HD11	2.01	0.42
35:BA:99:U:OP1	35:BA:102:G:OP1	2.38	0.42
35:BA:808:G:O2'	35:BA:1254:A:O2'	2.29	0.42
35:BA:1264:G:H2'	35:BA:1265:A:C8	2.55	0.42
35:BA:1285:G:H4'	48:BR:105:ARG:NH1	2.35	0.42
35:BA:1368:G:C6	35:BA:1369:G:N7	2.88	0.42
35:BA:1707:G:O4'	35:BA:1756:G:H1'	2.20	0.42
35:BA:207:A:H2'	35:BA:208:C:O4'	2.20	0.42
35:BA:2127:G:H1'	35:BA:2128:C:C4'	2.46	0.42
35:BA:2244:U:H1'	35:BA:2434:A:C8	2.55	0.42
35:BA:2538:C:C2'	35:BA:2539:C:C5'	2.98	0.42
35:BA:271(V):G:N3	35:BA:271(W):G:H1'	2.35	0.42
35:BA:2735:G:H2'	35:BA:2736:G:H8	1.83	0.42
35:BA:707:G:H3'	35:BA:708:C:C6	2.55	0.42
35:BA:71:A:C8	35:BA:71:A:H5'	2.54	0.42
31:B5:2:ALA:N	35:BA:747:U:C4	2.88	0.42
35:BA:770:G:C6	35:BA:771:G:N7	2.88	0.42
36:BB:10:C:C4	36:BB:11:C:C5	3.08	0.42
36:BB:16:G:O2'	36:BB:17:C:H5'	2.20	0.42
36:BB:29:A:OP2	49:BS:32:LEU:HB2	2.20	0.42
38:BD:133:LEU:CB	38:BD:173:VAL:HG11	2.49	0.42
35:BA:2228:G:OP2	38:BD:263:ARG:NH2	2.53	0.42
39:BE:143:ASN:N	39:BE:143:ASN:ND2	2.63	0.42
39:BE:2:LYS:NZ	39:BE:95:ILE:O	2.33	0.42
40:BF:150:GLY:HA2	40:BF:172:TRP:CD2	2.55	0.42
40:BF:160:ASN:HD21	40:BF:162:LEU:HB2	1.84	0.42
40:BF:20:LEU:O	40:BF:21:ALA:O	2.37	0.42
40:BF:82:ILE:C	40:BF:84:VAL:H	2.23	0.42
41:BG:104:GLU:O	41:BG:106:LEU:N	2.53	0.42
41:BG:106:LEU:O	41:BG:110:ALA:HB3	2.19	0.42
41:BG:17:PRO:C	41:BG:19:LEU:N	2.73	0.42
42:BH:16:SER:HB2	42:BH:27:LYS:O	2.20	0.42
42:BH:44:VAL:HG12	42:BH:45:VAL:HG23	2.02	0.42
43:BI:123:LEU:HD11	43:BI:143:SER:O	2.19	0.42
43:BI:55:ALA:O	43:BI:59:ALA:CB	2.68	0.42
44:BN:60:ILE:O	44:BN:61:ARG:C	2.58	0.42
46:BP:106:LEU:HB3	46:BP:107:LYS:H	1.62	0.42
35:BA:598:G:H5'	46:BP:15:ARG:CG	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:65:ARG:O	46:BP:66:GLY:C	2.57	0.42
48:BR:53:HIS:ND1	48:BR:53:HIS:O	2.52	0.42
48:BR:53:HIS:O	48:BR:56:LYS:HB2	2.20	0.42
48:BR:7:GLY:O	48:BR:8:ARG:O	2.38	0.42
49:BS:27:SER:O	49:BS:38:GLN:N	2.53	0.42
52:BV:82:ARG:CD	52:BV:82:ARG:C	2.88	0.42
53:BW:35:ILE:O	53:BW:36:LEU:C	2.57	0.42
53:BW:12:ILE:HD12	53:BW:42:ARG:NH1	2.34	0.42
35:BA:1341:U:H5'	54:BX:57:LEU:CD2	2.50	0.42
56:BZ:132:ASN:O	56:BZ:134:PRO:HD3	2.19	0.42
56:BZ:4:ARG:HA	56:BZ:59:LEU:HD23	2.02	0.42
1:CA:1197:G:C2'	1:CA:1198:G:H5'	2.50	0.42
1:CA:1357:A:N7	1:CA:1358:U:C4	2.87	0.42
1:CA:1357:A:N6	1:CA:1363(A):A:H2	2.16	0.42
1:CA:1400:C:H6	1:CA:1400:C:O5'	2.03	0.42
1:CA:178:C:O2'	1:CA:179:A:H5'	2.20	0.42
1:CA:226:G:O2'	1:CA:227:G:H5'	2.20	0.42
1:CA:349:A:C2'	1:CA:350:G:H5'	2.50	0.42
1:CA:460:G:C6	1:CA:470:C:H5''	2.55	0.42
1:CA:708:C:H2'	1:CA:709:G:H8	1.84	0.42
1:CA:722:A:O2'	1:CA:723:U:C6	2.72	0.42
2:CB:142:LEU:HA	2:CB:145:LEU:HB3	2.01	0.42
2:CB:69:LEU:HB2	2:CB:162:ILE:HG22	2.02	0.42
2:CB:197:VAL:HB	2:CB:200:ILE:HG12	2.01	0.42
3:CC:11:ARG:O	3:CC:13:GLY:N	2.53	0.42
4:CD:108:LEU:HD12	4:CD:108:LEU:N	2.34	0.42
4:CD:172:PRO:HD2	4:CD:173:TRP:HZ3	1.82	0.42
5:CE:139:LEU:H	5:CE:139:LEU:HG	1.53	0.42
5:CE:6:PHE:HB2	5:CE:34:VAL:HG12	2.01	0.42
5:CE:82:VAL:HG11	5:CE:134:ALA:O	2.20	0.42
8:CH:114:THR:CG2	8:CH:119:LEU:HD21	2.49	0.42
12:CL:47:LYS:HD3	12:CL:48:PRO:HD3	2.02	0.42
1:CA:1217:C:P	14:CN:5:ALA:HB1	2.60	0.42
1:CA:194:C:H4'	20:CT:65:LYS:HG3	2.01	0.42
25:CY:3:LEU:C	25:CY:5:GLU:N	2.71	0.42
28:D2:17:SER:HG	28:D2:18:PRO:HD3	1.83	0.42
29:D3:26:LEU:N	29:D3:26:LEU:HD23	2.35	0.42
32:D6:28:ARG:HA	32:D6:32:ASN:CB	2.50	0.42
32:D6:43:CYS:O	32:D6:44:ARG:O	2.37	0.42
33:D7:31:LEU:HD23	33:D7:42:LEU:HB3	2.01	0.42
35:DA:585:G:C4	35:DA:1251:C:N4	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:15:G:C4	35:DA:16:G:C8	3.08	0.42
35:DA:1619:G:N1	35:DA:1620:G:C5	2.88	0.42
35:DA:1831:G:H2'	35:DA:1832:C:H6	1.83	0.42
35:DA:1989:G:C2'	35:DA:1990:C:H5'	2.50	0.42
35:DA:2102:U:O4'	35:DA:2102:U:O2	2.35	0.42
35:DA:2061:G:O4'	35:DA:2503:A:C5	2.73	0.42
35:DA:2552:U:H2'	35:DA:2554:U:H5''	2.02	0.42
35:DA:225:A:O2'	35:DA:257:A:H4'	2.19	0.42
35:DA:2588:G:C6	35:DA:2589:A:C5	3.08	0.42
35:DA:2739:U:C2'	35:DA:2740:A:H5'	2.50	0.42
35:DA:2883:A:C5'	35:DA:2884:U:H5'	2.50	0.42
35:DA:2884:U:C2'	35:DA:2885:C:H5'	2.46	0.42
35:DA:340:A:H2'	35:DA:341:G:O4'	2.20	0.42
35:DA:607:U:H3	35:DA:621:A:H2	1.67	0.42
35:DA:723:G:C6	35:DA:724:U:C4	3.08	0.42
35:DA:7:G:H4'	44:DN:13:TRP:CZ2	2.54	0.42
35:DA:804:A:H2'	35:DA:806:C:C4	2.55	0.42
35:DA:915:C:H2'	35:DA:916:G:H8	1.85	0.42
36:DB:7:G:O5'	49:DS:29:PHE:HE2	2.02	0.42
39:DE:3:GLY:O	39:DE:4:ILE:CG2	2.66	0.42
39:DE:48:GLN:C	39:DE:49:LEU:HD22	2.40	0.42
41:DG:76:SER:HB3	41:DG:84:LYS:H	1.84	0.42
42:DH:46:GLU:CG	42:DH:51:ARG:HB2	2.50	0.42
44:DN:15:LEU:HD12	44:DN:136:GLU:CB	2.50	0.42
44:DN:63:THR:O	44:DN:64:GLY:C	2.58	0.42
46:DP:86:LYS:N	46:DP:117:GLU:O	2.52	0.42
46:DP:148:LEU:HD22	46:DP:148:LEU:C	2.41	0.42
47:DQ:140:ALA:HB2	56:DZ:99:TYR:CG	2.55	0.42
49:DS:38:GLN:CG	49:DS:39:ILE:N	2.80	0.42
49:DS:28:VAL:H	49:DS:89:ARG:CG	2.33	0.42
50:DT:84:GLN:HG3	50:DT:84:GLN:O	2.20	0.42
54:DX:61:GLY:N	54:DX:70:LEU:CD2	2.83	0.42
54:DX:88:LYS:CD	54:DX:88:LYS:N	2.83	0.42
56:DZ:150:LEU:O	56:DZ:171:ILE:CG1	2.68	0.42
56:DZ:17:ALA:HA	56:DZ:20:ARG:HB3	2.01	0.42
1:AA:1006:C:H2'	1:AA:1007:C:H6	1.85	0.42
1:AA:1299:A:N7	1:AA:1301:U:C2	2.88	0.42
1:AA:1349:A:H2'	1:AA:1350:A:O5'	2.20	0.42
1:AA:1428:A:C6	1:AA:1473:A:N1	2.87	0.42
1:AA:266:G:H5''	1:AA:268:C:N4	2.19	0.42
1:AA:533:A:C1'	1:AA:534:U:OP1	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:809:G:C2'	1:AA:810:C:O5'	2.68	0.42
2:AB:116:GLU:H	2:AB:116:GLU:HG2	1.55	0.42
2:AB:213:LEU:CD2	2:AB:214:ILE:N	2.83	0.42
3:AC:87:LEU:CB	3:AC:101:LEU:HD11	2.48	0.42
4:AD:150:GLU:CA	4:AD:153:ARG:HG3	2.50	0.42
5:AE:144:THR:C	5:AE:148:VAL:HG23	2.40	0.42
6:AF:37:VAL:HG12	6:AF:38:GLU:N	2.35	0.42
8:AH:111:ILE:O	8:AH:112:LEU:HB3	2.20	0.42
8:AH:5:PRO:O	8:AH:8:ASP:HB3	2.20	0.42
11:AK:86:GLY:H	11:AK:112:THR:CB	2.33	0.42
13:AM:108:ARG:O	13:AM:109:THR:C	2.58	0.42
13:AM:4:ILE:CG2	13:AM:5:ALA:H	2.33	0.42
1:AA:1219:U:P	14:AN:19:ARG:HH22	2.42	0.42
16:AP:53:VAL:O	16:AP:54:GLU:C	2.58	0.42
6:AF:97:PHE:O	18:AR:30:ASP:HA	2.20	0.42
20:AT:22:ARG:O	20:AT:26:ASN:ND2	2.53	0.42
20:AT:27:LYS:O	20:AT:30:LYS:HB3	2.20	0.42
21:AU:17:THR:O	21:AU:22:ARG:NH1	2.52	0.42
21:AU:24:ARG:H	21:AU:24:ARG:HD2	1.85	0.42
23:AW:2:G:C6	23:AW:73:A:C2	3.08	0.42
25:AY:29:ARG:H	25:AY:29:ARG:HG2	1.43	0.42
32:B6:23:THR:HG21	35:BA:2419:U:H5'	2.02	0.42
33:B7:1:MET:O	35:BA:1620:G:O4'	2.38	0.42
35:BA:122:G:H1	35:BA:129:C:N4	2.17	0.42
35:BA:1353:A:H4'	38:BD:38:LYS:HZ1	1.80	0.42
35:BA:1313:U:H2'	35:BA:1610:A:C2	2.54	0.42
35:BA:1623:G:H2'	35:BA:1624:G:C8	2.51	0.42
35:BA:1810:A:C2'	35:BA:1811:G:H5'	2.50	0.42
35:BA:205:G:O2'	35:BA:206:U:OP2	2.36	0.42
35:BA:2200:C:C6	35:BA:2200:C:H5''	2.54	0.42
35:BA:415:A:N1	35:BA:2409:G:C6	2.88	0.42
34:B8:35:GLN:HG2	35:BA:2420:C:P	2.59	0.42
35:BA:2697:G:H2'	35:BA:2698:U:O4'	2.20	0.42
35:BA:2881:C:C4	35:BA:2882:A:N7	2.88	0.42
35:BA:341:G:H2'	35:BA:342:G:C8	2.54	0.42
35:BA:261:G:H1'	35:BA:609:A:H2	1.84	0.42
35:BA:648:G:O4'	35:BA:2351:G:H5''	2.20	0.42
35:BA:656:G:C6	35:BA:657:U:C4	3.08	0.42
35:BA:966:G:H2'	35:BA:967:C:C6	2.55	0.42
35:BA:996:A:H2'	35:BA:997:G:C8	2.51	0.42
36:BB:79:C:H42	36:BB:98:G:H1	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BF:160:ASN:CB	40:BF:163:VAL:HG23	2.48	0.42
30:B4:6:HIS:CB	41:BG:67:LYS:CE	2.98	0.42
42:BH:149:ARG:HE	42:BH:154:PRO:HD3	1.83	0.42
43:BI:114:LEU:O	43:BI:115:ALA:CB	2.67	0.42
44:BN:63:THR:O	44:BN:64:GLY:C	2.57	0.42
45:BO:87:ILE:HG21	45:BO:91:LEU:HD13	2.02	0.42
46:BP:75:ILE:O	46:BP:76:LYS:C	2.57	0.42
47:BQ:70:PRO:CA	47:BQ:95:ALA:HB2	2.50	0.42
35:BA:1653:G:O6	48:BR:11:ASN:HB2	2.20	0.42
48:BR:32:GLY:C	48:BR:33:ARG:HG3	2.40	0.42
48:BR:26:LYS:HZ3	48:BR:71:GLN:HB3	1.83	0.42
53:BW:64:MET:HG2	53:BW:109:GLU:OE2	2.20	0.42
53:BW:57:ASN:HA	53:BW:57:ASN:HD22	1.63	0.42
53:BW:67:ASP:N	53:BW:69:LEU:HD11	2.33	0.42
35:BA:143:G:C1'	54:BX:38:GLU:HG3	2.50	0.42
55:BY:88:LYS:CD	55:BY:88:LYS:N	2.79	0.42
56:BZ:103:ARG:HB2	56:BZ:138:GLU:CA	2.37	0.42
56:BZ:5:LEU:HB2	56:BZ:59:LEU:HD22	2.02	0.42
56:BZ:61:LEU:HD12	56:BZ:65:GLN:HB3	2.01	0.42
56:BZ:63:ASP:C	56:BZ:65:GLN:N	2.73	0.42
56:BZ:63:ASP:O	56:BZ:65:GLN:N	2.52	0.42
1:CA:1168:A:C2	1:CA:1169:A:C4	3.07	0.42
1:CA:1305:G:C2	1:CA:1331:G:N3	2.87	0.42
1:CA:1237:C:C4'	1:CA:1334:G:N2	2.83	0.42
1:CA:1342:C:H2'	1:CA:1343:G:H8	1.84	0.42
1:CA:1355:G:O2'	1:CA:1356:G:H5'	2.20	0.42
1:CA:1365:G:H2'	1:CA:1366:C:C5'	2.50	0.42
1:CA:1064:G:OP2	1:CA:1386:G:H4'	2.20	0.42
1:CA:925:G:H1	1:CA:1391:U:H3	1.66	0.42
1:CA:1396:A:N3	5:CE:19:MET:HG3	2.35	0.42
1:CA:1517:G:H2'	1:CA:1518:A:O5'	2.20	0.42
1:CA:545:C:OP1	4:CD:61:LYS:NZ	2.52	0.42
1:CA:552:U:H2'	1:CA:553:A:H8	1.84	0.42
1:CA:608:A:H2'	1:CA:609:A:O4'	2.19	0.42
1:CA:696:A:C4	1:CA:697:U:C5	3.08	0.42
4:CD:150:GLU:CA	4:CD:153:ARG:HG3	2.50	0.42
1:CA:15:G:H1'	5:CE:19:MET:HG2	2.02	0.42
6:CF:7:ASN:O	6:CF:8:ILE:CG1	2.60	0.42
7:CG:47:CYS:C	7:CG:58:PRO:HG3	2.40	0.42
7:CG:66:VAL:O	7:CG:69:VAL:N	2.40	0.42
8:CH:35:ILE:O	8:CH:36:LEU:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:4:ASP:OD2	8:CH:7:ALA:CB	2.67	0.42
1:CA:1525:G:P	11:CK:120:ARG:HH22	2.43	0.42
12:CL:55:VAL:C	12:CL:70:ILE:HD11	2.40	0.42
13:CM:28:ALA:O	13:CM:32:GLU:HB2	2.20	0.42
15:CO:12:ILE:HG23	15:CO:27:VAL:HG11	2.00	0.42
15:CO:30:ALA:HA	15:CO:85:LEU:HD11	2.00	0.42
16:CP:19:ILE:HB	16:CP:37:GLY:C	2.40	0.42
27:D1:25:LYS:HB2	27:D1:37:ILE:CG2	2.50	0.42
27:D1:58:ILE:CG2	27:D1:59:THR:N	2.75	0.42
29:D3:11:SER:OG	29:D3:13:ILE:HG12	2.19	0.42
29:D3:16:PRO:CB	29:D3:18:ASP:OD1	2.68	0.42
31:D5:15:ARG:HG3	31:D5:15:ARG:NH1	2.35	0.42
32:D6:15:GLU:O	32:D6:16:CYS:O	2.38	0.42
32:D6:12:GLU:O	32:D6:51:GLU:O	2.37	0.42
34:D8:7:HIS:HD2	46:DP:50:ARG:CZ	2.32	0.42
35:DA:127:A:H5'	35:DA:128:C:C1'	2.50	0.42
35:DA:1300:U:O2'	35:DA:1301:A:OP2	2.35	0.42
35:DA:1500:G:C6	35:DA:1501:C:N3	2.87	0.42
35:DA:1766:U:H2'	35:DA:1767:C:H6	1.83	0.42
35:DA:1790:C:H2'	35:DA:1791:A:C5	2.54	0.42
35:DA:203:C:H2'	35:DA:204:A:H8	1.84	0.42
35:DA:234:C:O5'	35:DA:234:C:H6	2.02	0.42
35:DA:2625:G:H2'	35:DA:2626:C:O4'	2.20	0.42
35:DA:2687:U:C2'	35:DA:2688:U:H5'	2.50	0.42
35:DA:2881:C:N3	35:DA:2882:A:N7	2.68	0.42
35:DA:29:U:H2'	35:DA:30:G:C8	2.55	0.42
35:DA:261:G:H1'	35:DA:609:A:H2	1.85	0.42
35:DA:688:U:H5'	35:DA:1780:A:C2	2.55	0.42
35:DA:58:G:N2	35:DA:70:G:C4	2.88	0.42
35:DA:806:C:O2	35:DA:2444:G:O2'	2.37	0.42
35:DA:855:G:C6	35:DA:856:C:C4	3.08	0.42
36:DB:40:U:C2'	36:DB:41:U:OP1	2.68	0.42
38:DD:49:ILE:HG13	38:DD:49:ILE:O	2.20	0.42
39:DE:31:CYS:HA	39:DE:50:GLY:O	2.19	0.42
39:DE:53:PRO:O	39:DE:54:GLN:C	2.58	0.42
39:DE:72:VAL:O	39:DE:73:GLU:C	2.57	0.42
40:DF:153:SER:O	40:DF:190:GLU:HB2	2.20	0.42
41:DG:101:ILE:O	41:DG:104:GLU:HB2	2.19	0.42
41:DG:115:ARG:HH22	41:DG:136:ARG:HD2	1.85	0.42
43:DI:71:ILE:HG13	43:DI:72:LEU:CD2	2.50	0.42
43:DI:82:ARG:HG3	43:DI:82:ARG:NH1	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:64:ARG:HB2	45:DO:64:ARG:HH11	1.85	0.42
46:DP:16:ARG:O	46:DP:16:ARG:HD3	2.19	0.42
48:DR:29:LEU:HG	48:DR:79:LEU:CD2	2.49	0.42
48:DR:26:LYS:HZ3	48:DR:71:GLN:HB3	1.84	0.42
50:DT:17:THR:O	50:DT:18:ASP:HB3	2.20	0.42
52:DV:39:LEU:HD11	52:DV:53:GLU:H	1.83	0.42
52:DV:61:VAL:O	52:DV:62:LEU:HD23	2.19	0.42
52:DV:25:LEU:N	52:DV:94:LEU:HD11	2.34	0.42
53:DW:12:ILE:HD12	53:DW:42:ARG:HH11	1.84	0.42
28:D2:23:LYS:N	54:DX:5:TYR:CE1	2.88	0.42
54:DX:83:VAL:C	54:DX:85:PRO:CD	2.88	0.42
55:DY:81:LYS:HD3	55:DY:97:ARG:C	2.39	0.42
56:DZ:141:VAL:O	56:DZ:142:SER:O	2.36	0.42
1:AA:1054:C:OP1	1:AA:1197:G:OP2	2.38	0.41
1:AA:1126:U:H2'	1:AA:1127:G:O4'	2.20	0.41
1:AA:1164:G:C2'	1:AA:1165:C:H5'	2.50	0.41
1:AA:1197:G:C2'	1:AA:1198:G:H5'	2.50	0.41
1:AA:1249:C:H5'	1:AA:1249:C:H6	1.85	0.41
1:AA:1254:C:H2'	1:AA:1255:G:H8	1.85	0.41
1:AA:1298:C:O4'	1:AA:1299:A:C4	2.73	0.41
1:AA:296:U:O2'	1:AA:297:G:H5'	2.20	0.41
1:AA:30:U:O2'	1:AA:31:G:OP1	2.29	0.41
1:AA:541:G:H2'	1:AA:542:G:H8	1.84	0.41
1:AA:768:A:N3	1:AA:1512:U:O2'	2.53	0.41
1:AA:73:G:C2	1:AA:97:G:N1	2.88	0.41
2:AB:153:ARG:O	2:AB:154:LEU:O	2.38	0.41
2:AB:211:ILE:O	2:AB:215:LEU:CB	2.68	0.41
2:AB:222:ILE:CG2	2:AB:223:ILE:N	2.83	0.41
2:AB:23:ARG:O	2:AB:23:ARG:HG2	2.20	0.41
2:AB:70:PHE:CD1	2:AB:163:PHE:HB3	2.54	0.41
3:AC:59:ARG:HA	3:AC:63:ASN:O	2.20	0.41
4:AD:202:LEU:O	4:AD:204:ILE:N	2.53	0.41
5:AE:144:THR:O	5:AE:146:ALA:N	2.53	0.41
6:AF:14:LEU:HD11	6:AF:19:LEU:HB2	1.98	0.41
7:AG:111:ARG:HB3	7:AG:112:PRO:HD2	2.02	0.41
7:AG:39:ALA:O	7:AG:41:ARG:N	2.53	0.41
7:AG:74:GLU:CG	7:AG:75:VAL:N	2.83	0.41
7:AG:74:GLU:HG2	7:AG:75:VAL:N	2.35	0.41
8:AH:95:VAL:HG22	8:AH:131:GLY:O	2.20	0.41
9:AI:92:TYR:N	9:AI:92:TYR:CD1	2.88	0.41
10:AJ:50:ILE:HA	10:AJ:60:ARG:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AJ:51:ARG:HG3	10:AJ:60:ARG:C	2.41	0.41
10:AJ:48:THR:HG1	10:AJ:62:HIS:CG	2.38	0.41
13:AM:40:ASN:HD22	13:AM:43:THR:HG23	1.83	0.41
13:AM:78:ILE:HA	13:AM:81:LEU:HD12	2.01	0.41
16:AP:55:ARG:O	16:AP:58:TYR:HB3	2.19	0.41
1:AA:1320:C:H42	19:AS:36:ARG:HG3	1.84	0.41
20:AT:58:LYS:HE3	20:AT:62:LEU:HD11	2.01	0.41
25:AY:18:LEU:CD2	25:AY:171:LYS:HB3	2.49	0.41
25:AY:52:LEU:O	25:AY:54:GLN:N	2.53	0.41
25:AY:79:ILE:HG23	25:AY:83:ILE:HD11	2.01	0.41
28:B2:18:PRO:O	28:B2:22:GLU:N	2.53	0.41
30:B4:13:ARG:O	30:B4:15:ILE:N	2.42	0.41
31:B5:40:LYS:NZ	31:B5:50:GLY:HA2	2.35	0.41
35:BA:83:G:C4	35:BA:102:G:N2	2.88	0.41
35:BA:1433:U:O2	35:BA:1561:G:C2	2.73	0.41
35:BA:1619:G:N1	35:BA:1620:G:C5	2.88	0.41
35:BA:1635:G:H2'	35:BA:1636:C:C6	2.55	0.41
35:BA:1775:U:C2'	35:BA:1776:G:O5'	2.68	0.41
35:BA:1935:G:H3'	35:BA:1962:C:N4	2.32	0.41
35:BA:2032:G:N2	35:BA:2572:A:C8	2.88	0.41
26:B0:16:SER:OG	35:BA:2261:C:H3'	2.20	0.41
35:BA:2382:G:C3'	35:BA:2383:G:H5'	2.50	0.41
35:BA:2392:A:C2	35:BA:2429:G:C4	3.08	0.41
35:BA:2476:A:C3'	35:BA:2477:C:H5''	2.50	0.41
35:BA:2531:A:H2	35:BA:2658:C:O2	2.03	0.41
35:BA:2572:A:P	39:BE:144:ARG:HB2	2.60	0.41
35:BA:2618:G:H2'	35:BA:2619:C:H6	1.84	0.41
35:BA:2685:G:OP1	50:BT:51:ARG:NH2	2.53	0.41
35:BA:2880:C:H4'	48:BR:90:ARG:NH1	2.35	0.41
35:BA:14:A:C6	35:BA:526:A:C2	3.07	0.41
35:BA:566:U:H4'	35:BA:809:G:OP2	2.20	0.41
35:BA:577:G:N1	35:BA:578:A:C6	2.88	0.41
35:BA:8:A:OP1	44:BN:51:PHE:HE2	2.03	0.41
35:BA:904:C:H2'	35:BA:905:U:H5'	2.01	0.41
35:BA:986:C:C2'	35:BA:987:G:H5'	2.50	0.41
36:BB:82:G:O2'	36:BB:83:G:H5'	2.19	0.41
37:BC:18:LYS:O	37:BC:20:TYR:N	2.53	0.41
37:BC:42:GLU:O	37:BC:212:VAL:HA	2.19	0.41
37:BC:51:PRO:HB3	37:BC:204:ALA:CB	2.50	0.41
38:BD:140:THR:O	38:BD:165:ILE:CD1	2.68	0.41
39:BE:35:GLN:NE2	39:BE:37:ARG:NH2	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:3:GLY:O	39:BE:4:ILE:HB	2.20	0.41
41:BG:140:ILE:HD12	41:BG:140:ILE:C	2.40	0.41
41:BG:174:GLU:HG3	41:BG:182:LYS:HZ1	1.84	0.41
42:BH:127:GLU:HB2	42:BH:130:ARG:HB3	2.00	0.41
42:BH:122:THR:O	42:BH:133:VAL:HG22	2.20	0.41
42:BH:94:TYR:CZ	42:BH:160:LYS:HD2	2.54	0.41
43:BI:114:LEU:O	43:BI:129:THR:O	2.37	0.41
43:BI:37:VAL:HG12	43:BI:38:LEU:N	2.34	0.41
35:BA:943:U:OP2	46:BP:38:GLN:CG	2.68	0.41
47:BQ:120:ILE:HA	47:BQ:123:HIS:HD2	1.84	0.41
48:BR:14:SER:O	48:BR:15:SER:C	2.59	0.41
48:BR:81:ASP:O	48:BR:82:GLU:HB2	2.20	0.41
49:BS:36:TYR:O	49:BS:37:ALA:HB2	2.20	0.41
49:BS:68:GLN:O	49:BS:71:ARG:HB2	2.20	0.41
50:BT:101:PHE:C	50:BT:101:PHE:CD2	2.93	0.41
45:BO:104:ARG:NH2	50:BT:33:LYS:CE	2.83	0.41
54:BX:36:LYS:CD	54:BX:36:LYS:O	2.61	0.41
54:BX:78:LYS:C	54:BX:78:LYS:HD3	2.40	0.41
56:BZ:109:ALA:O	56:BZ:110:GLY:O	2.37	0.41
1:CA:1007:C:H2'	1:CA:1008:C:C5	2.54	0.41
1:CA:1071:C:H2'	1:CA:1072:G:H8	1.84	0.41
1:CA:1084:G:C5	1:CA:1085:U:C4	3.07	0.41
1:CA:1187:G:H2'	1:CA:1187:G:N3	2.35	0.41
1:CA:1248:A:H2'	1:CA:1249:C:C5'	2.49	0.41
1:CA:1284:C:H3'	1:CA:1285:A:H8	1.84	0.41
1:CA:1510:U:C2	1:CA:1526:G:N2	2.88	0.41
1:CA:540:G:H2'	1:CA:541:G:O4'	2.20	0.41
1:CA:663:A:H2'	1:CA:664:G:C8	2.54	0.41
1:CA:686:U:H1'	1:CA:687:A:C8	2.55	0.41
1:CA:762:C:H2'	1:CA:763:G:H8	1.84	0.41
1:CA:799:G:C6	1:CA:800:G:C4	3.08	0.41
1:CA:880:C:H2'	1:CA:881:G:C8	2.55	0.41
1:CA:983:A:N1	1:CA:1222:G:N2	2.67	0.41
4:CD:162:LEU:O	4:CD:165:MET:HB2	2.19	0.41
5:CE:62:ALA:O	5:CE:63:ARG:C	2.58	0.41
6:CF:62:TRP:C	6:CF:63:TYR:CD2	2.93	0.41
7:CG:78:ARG:HG3	7:CG:79:ARG:N	2.35	0.41
8:CH:134:ILE:HG22	8:CH:135:CYS:SG	2.59	0.41
11:CK:20:TYR:O	11:CK:31:THR:N	2.51	0.41
12:CL:117:ARG:HD2	12:CL:122:THR:HB	2.02	0.41
13:CM:83:ASP:OD2	13:CM:84:ILE:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CO:8:LYS:HG2	15:CO:12:ILE:HD11	2.02	0.41
1:CA:657:G:N2	15:CO:22:THR:OG1	2.52	0.41
15:CO:51:HIS:O	15:CO:52:SER:C	2.57	0.41
17:CQ:17:LYS:HA	17:CQ:46:ASP:O	2.19	0.41
17:CQ:45:HIS:HE2	17:CQ:47:PRO:HB3	1.85	0.41
22:CV:34:A:H2'	22:CV:35:A:C8	2.55	0.41
23:CW:72:C:H2'	23:CW:73:A:O4'	2.20	0.41
27:D1:27:GLU:OE1	27:D1:32:LYS:HB2	2.20	0.41
27:D1:57:GLU:HA	27:D1:57:GLU:OE1	2.20	0.41
27:D1:84:GLY:O	27:D1:85:LEU:C	2.58	0.41
29:D3:6:VAL:HG12	29:D3:56:VAL:HA	2.02	0.41
31:D5:40:LYS:HZ3	31:D5:45:VAL:CA	2.29	0.41
32:D6:12:GLU:OE1	32:D6:12:GLU:N	2.53	0.41
35:DA:1039:G:N1	35:DA:1117:G:C2	2.88	0.41
35:DA:1766:U:H2'	35:DA:1767:C:C6	2.54	0.41
35:DA:1792:G:OP1	38:DD:206:LEU:HB2	2.20	0.41
35:DA:1810:A:H2'	35:DA:1811:G:O4'	2.20	0.41
35:DA:2171:A:O2'	35:DA:2172:U:H6	2.03	0.41
35:DA:2555:U:H2'	35:DA:2556:C:C5'	2.49	0.41
35:DA:324:A:N6	35:DA:338:G:O2'	2.51	0.41
35:DA:373:U:O2	35:DA:423:A:H2	2.03	0.41
35:DA:426:C:H2'	35:DA:427:U:C5'	2.50	0.41
35:DA:479:A:N6	35:DA:503:A:H61	2.17	0.41
35:DA:531:C:C5	35:DA:2035:G:C2	3.08	0.41
35:DA:562:U:C2'	35:DA:563:G:OP2	2.67	0.41
38:DD:130:ALA:HB2	38:DD:192:THR:CA	2.50	0.41
38:DD:36:PRO:CA	38:DD:62:TYR:O	2.68	0.41
38:DD:35:LYS:CG	38:DD:64:ILE:N	2.70	0.41
39:DE:82:ARG:CG	39:DE:83:ASP:H	2.26	0.41
40:DF:110:LEU:HD11	40:DF:181:LEU:O	2.20	0.41
40:DF:53:THR:HG23	40:DF:56:GLU:H	1.85	0.41
40:DF:80:ALA:O	40:DF:83:PHE:HB2	2.20	0.41
41:DG:102:PHE:O	41:DG:103:LEU:C	2.58	0.41
41:DG:105:LYS:HB3	41:DG:142:PRO:CG	2.50	0.41
41:DG:48:GLU:O	41:DG:49:ASP:HB2	2.20	0.41
41:DG:59:GLU:CD	41:DG:60:LEU:CD2	2.89	0.41
41:DG:4:ASP:CA	41:DG:8:LYS:HD3	2.42	0.41
42:DH:74:ASN:O	42:DH:76:VAL:N	2.52	0.41
42:DH:92:ILE:CG2	42:DH:93:GLY:N	2.72	0.41
44:DN:51:PHE:O	44:DN:52:VAL:C	2.58	0.41
44:DN:60:ILE:HD13	44:DN:99:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DO:104:ARG:HH21	50:DT:33:LYS:CE	2.29	0.41
45:DO:13:ASN:ND2	45:DO:97:ARG:N	2.68	0.41
35:DA:598:G:H5'	46:DP:15:ARG:CG	2.49	0.41
46:DP:18:ARG:HB3	46:DP:18:ARG:HE	1.53	0.41
47:DQ:24:GLY:CA	47:DQ:101:ARG:HA	2.50	0.41
47:DQ:61:GLY:O	47:DQ:62:GLY:O	2.38	0.41
47:DQ:14:ARG:O	47:DQ:72:LYS:HE2	2.19	0.41
48:DR:2:ARG:CD	48:DR:2:ARG:C	2.88	0.41
48:DR:32:GLY:O	48:DR:116:LEU:N	2.52	0.41
49:DS:54:LEU:CD1	49:DS:58:LEU:O	2.63	0.41
50:DT:101:PHE:C	50:DT:101:PHE:CD2	2.93	0.41
50:DT:27:THR:HA	50:DT:88:ILE:N	2.30	0.41
50:DT:28:VAL:HG23	50:DT:47:GLY:O	2.19	0.41
50:DT:53:ARG:CG	50:DT:53:ARG:O	2.68	0.41
50:DT:8:LYS:O	50:DT:8:LYS:HG2	2.20	0.41
52:DV:99:ILE:HG22	52:DV:100:ARG:N	2.35	0.41
52:DV:39:LEU:HD22	52:DV:53:GLU:O	2.20	0.41
35:DA:1162:G:N2	52:DV:91:TYR:HE1	2.18	0.41
53:DW:71:VAL:HG12	53:DW:71:VAL:O	2.19	0.41
56:DZ:110:GLY:C	56:DZ:111:VAL:HG12	2.40	0.41
56:DZ:19:ARG:CZ	56:DZ:84:GLU:OE2	2.68	0.41
1:AA:1050:G:H2'	1:AA:1051:C:H6	1.85	0.41
1:AA:1059:C:O2	10:AJ:53:PRO:HG3	2.20	0.41
1:AA:1223:C:H3'	1:AA:1224:G:H5''	2.02	0.41
1:AA:1511:G:O5'	1:AA:1511:G:H8	2.04	0.41
1:AA:298:A:H2'	1:AA:299:G:O4'	2.20	0.41
1:AA:452:A:C4	1:AA:453:A:C8	3.09	0.41
1:AA:44:G:H2'	1:AA:45:U:O4'	2.20	0.41
1:AA:947:G:H2'	1:AA:948:C:C6	2.54	0.41
1:AA:983:A:N1	1:AA:1222:G:N2	2.67	0.41
1:AA:992:U:O2'	1:AA:993:G:P	2.78	0.41
2:AB:212:GLN:O	2:AB:213:LEU:C	2.58	0.41
2:AB:36:ARG:NE	2:AB:36:ARG:HA	2.35	0.41
4:AD:100:ARG:HG2	4:AD:100:ARG:HH11	1.84	0.41
7:AG:24:THR:HA	7:AG:27:ILE:HD13	2.01	0.41
7:AG:65:ALA:HB2	7:AG:124:LEU:O	2.20	0.41
9:AI:70:LYS:O	9:AI:73:GLN:HB2	2.20	0.41
1:AA:973:G:P	10:AJ:57:LYS:NZ	2.93	0.41
11:AK:114:VAL:O	11:AK:114:VAL:HG13	2.20	0.41
11:AK:84:VAL:HG11	11:AK:95:ILE:CD1	2.50	0.41
12:AL:90:VAL:C	12:AL:92:ASP:N	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AM:91:ARG:NH2	13:AM:97:PRO:O	2.53	0.41
15:AO:67:LEU:HA	15:AO:67:LEU:HD23	1.80	0.41
12:AL:10:LEU:HB3	17:AQ:32:TYR:CE1	2.55	0.41
18:AR:22:VAL:HG13	18:AR:25:THR:HB	2.01	0.41
19:AS:45:VAL:C	19:AS:47:HIS:H	2.22	0.41
22:AV:31:U:H6	22:AV:31:U:O5'	2.03	0.41
23:AW:17:C:O5'	23:AW:62:C:H5'	2.20	0.41
23:AW:6:G:C2	23:AW:69:C:O2	2.73	0.41
26:B0:36:ILE:HD11	35:BA:2355:C:O4'	2.20	0.41
26:B0:70:GLN:HB3	26:B0:78:TYR:HB2	2.02	0.41
27:B1:13:ILE:CB	27:B1:63:ALA:HB2	2.46	0.41
27:B1:70:VAL:O	27:B1:72:GLU:N	2.53	0.41
31:B5:15:ARG:HA	31:B5:18:ALA:HB3	2.01	0.41
32:B6:12:GLU:O	32:B6:51:GLU:O	2.37	0.41
35:BA:998:C:N4	35:BA:1158:C:N4	2.67	0.41
35:BA:1176:G:C1'	35:BA:1177:A:OP1	2.64	0.41
35:BA:1417:C:H2'	35:BA:1418:G:C5'	2.50	0.41
35:BA:2071:A:H2'	35:BA:2072:G:C8	2.54	0.41
35:BA:2219:G:C2'	35:BA:2220:G:H5'	2.50	0.41
35:BA:2297:C:H2'	35:BA:2298:A:C5'	2.36	0.41
35:BA:2313:C:H2'	35:BA:2314:C:C6	2.55	0.41
35:BA:2331:G:C6	35:BA:2332:U:C4	3.08	0.41
35:BA:2627:G:H21	35:BA:2781:A:H2	1.65	0.41
35:BA:2732:G:H2'	35:BA:2733:A:H5'	1.97	0.41
35:BA:2765:A:H2	35:BA:2766:G:O4'	2.03	0.41
35:BA:2801(A):A:H4'	35:BA:2802:G:C5'	2.48	0.41
35:BA:2818:G:O2'	35:BA:2819:G:H5'	2.20	0.41
35:BA:42:G:C2	35:BA:437:G:C2	3.08	0.41
35:BA:28:A:H1'	35:BA:513:A:N1	2.35	0.41
35:BA:863:A:H4'	36:BB:101:G:C2	2.55	0.41
36:BB:16:G:H2'	36:BB:17:C:H6	1.84	0.41
36:BB:71:C:H2'	36:BB:72:G:O4'	2.20	0.41
36:BB:81:G:H5'	36:BB:82:G:OP2	2.20	0.41
37:BC:86:ALA:HB1	37:BC:94:VAL:HG11	2.02	0.41
38:BD:226:MET:CB	38:BD:230:ASP:HB2	2.41	0.41
38:BD:94:LEU:C	38:BD:94:LEU:HD13	2.40	0.41
40:BF:123:LEU:HD13	40:BF:192:LEU:HD22	2.02	0.41
42:BH:61:HIS:C	42:BH:63:SER:H	2.22	0.41
43:BI:60:GLU:C	43:BI:62:LYS:H	2.22	0.41
44:BN:113:GLY:O	44:BN:114:ARG:C	2.57	0.41
44:BN:79:PRO:HG2	44:BN:80:GLY:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BP:148:LEU:HD22	46:BP:148:LEU:C	2.40	0.41
47:BQ:68:ILE:HG22	47:BQ:101:ARG:HE	1.85	0.41
47:BQ:16:ARG:HH11	47:BQ:16:ARG:CB	2.33	0.41
50:BT:32:TYR:HD2	50:BT:81:PRO:HB2	1.85	0.41
52:BV:99:ILE:HG22	52:BV:100:ARG:N	2.34	0.41
53:BW:71:VAL:HA	53:BW:107:LEU:HD12	2.01	0.41
54:BX:76:ARG:O	54:BX:77:LYS:CB	2.66	0.41
56:BZ:177:PRO:O	56:BZ:178:GLU:CG	2.61	0.41
1:CA:1397:C:N3	24:CX:22:U:H6	2.17	0.41
1:CA:1422:G:C2	1:CA:1423:G:N7	2.88	0.41
1:CA:1445:C:O2'	1:CA:1446:U:H5'	2.20	0.41
1:CA:1458:G:C4	1:CA:1459:C:C5	3.08	0.41
1:CA:474:G:H2'	1:CA:475:G:H8	1.86	0.41
1:CA:825:G:O2'	1:CA:826:C:H5'	2.20	0.41
3:CC:134:ILE:HG23	3:CC:151:VAL:CG1	2.49	0.41
3:CC:150:LYS:HA	3:CC:168:ALA:O	2.20	0.41
4:CD:39:PRO:HB3	4:CD:40:PRO:HD2	2.01	0.41
4:CD:60:GLU:O	4:CD:63:LYS:N	2.53	0.41
6:CF:14:LEU:HD11	6:CF:19:LEU:HB2	2.01	0.41
7:CG:60:LYS:HD2	7:CG:63:LYS:HG2	2.02	0.41
7:CG:66:VAL:HB	7:CG:67:GLU:OE2	2.20	0.41
7:CG:74:GLU:HG2	7:CG:75:VAL:N	2.35	0.41
8:CH:36:LEU:O	8:CH:37:ARG:C	2.57	0.41
8:CH:26:VAL:HG12	8:CH:59:LEU:HB2	2.01	0.41
1:CA:1347:G:C2	9:CI:107:ARG:NH2	2.88	0.41
1:CA:1249:C:H4'	9:CI:36:TYR:OH	2.19	0.41
11:CK:116:HIS:O	11:CK:117:ASN:CB	2.67	0.41
11:CK:29:ILE:HD12	11:CK:29:ILE:O	2.19	0.41
12:CL:70:ILE:HG13	12:CL:100:ILE:CD1	2.49	0.41
13:CM:43:THR:HB	13:CM:44:ARG:H	1.61	0.41
13:CM:4:ILE:HG22	13:CM:5:ALA:N	2.35	0.41
14:CN:41:ARG:CG	14:CN:42:ILE:H	2.33	0.41
15:CO:67:LEU:HA	15:CO:67:LEU:HD23	1.79	0.41
16:CP:74:LEU:O	16:CP:77:ALA:HB3	2.20	0.41
17:CQ:12:SER:HA	17:CQ:14:LYS:HZ1	1.86	0.41
1:CA:1457:G:P	20:CT:39:LYS:HZ2	2.43	0.41
25:CY:131:ASN:O	25:CY:133:ARG:N	2.53	0.41
25:CY:16:LYS:C	25:CY:18:LEU:H	2.23	0.41
27:D1:76:ARG:CA	27:D1:76:ARG:NE	2.82	0.41
28:D2:29:LYS:O	28:D2:32:LEU:N	2.52	0.41
35:DA:1389:G:H2'	35:DA:1390:U:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1564:C:H2'	35:DA:1565:C:C6	2.55	0.41
35:DA:1577:C:H2'	35:DA:1578:U:H6	1.77	0.41
35:DA:1271:G:N2	35:DA:1617:C:H4'	2.35	0.41
35:DA:1301:A:H2	35:DA:1626:G:H21	1.68	0.41
35:DA:1655:A:H4'	39:DE:115:GLY:N	2.34	0.41
35:DA:1688:U:O2	35:DA:1700:A:H8	2.04	0.41
35:DA:1722:A:H2	35:DA:1740:G:H2'	1.86	0.41
35:DA:1707:G:C4	35:DA:1756:G:C2	3.08	0.41
35:DA:1778:U:C6	35:DA:1784:A:C6	3.08	0.41
35:DA:1783:A:OP1	35:DA:1784:A:OP2	2.38	0.41
35:DA:1790:C:C2'	35:DA:1791:A:C8	3.03	0.41
35:DA:1666:G:N1	35:DA:1995:U:O4	2.53	0.41
35:DA:2163:C:O2'	35:DA:2164:C:P	2.78	0.41
35:DA:225:A:N6	35:DA:419:C:H4'	2.36	0.41
35:DA:2343:C:O2'	35:DA:2373:G:H4'	2.19	0.41
35:DA:2597:G:C5	35:DA:2598:A:N6	2.88	0.41
35:DA:2633:G:C2	35:DA:2634:G:C4	3.08	0.41
35:DA:447:A:N3	35:DA:473:G:C8	2.88	0.41
35:DA:614:U:C4'	35:DA:614(C):A:H62	2.33	0.41
36:DB:43:C:H3'	36:DB:44:G:H5'	2.01	0.41
36:DB:64:C:H2'	36:DB:65:C:C6	2.55	0.41
38:DD:211:ARG:C	38:DD:213:ARG:N	2.73	0.41
40:DF:182:ASN:OD1	40:DF:185:ASP:N	2.53	0.41
40:DF:199:TRP:O	40:DF:202:PHE:HB3	2.21	0.41
41:DG:19:LEU:HA	41:DG:22:ARG:HB3	2.02	0.41
42:DH:121:ILE:HG22	42:DH:133:VAL:HG11	2.02	0.41
42:DH:68:THR:C	42:DH:70:THR:N	2.73	0.41
44:DN:116:LEU:O	44:DN:119:ARG:HB2	2.20	0.41
44:DN:62:VAL:O	44:DN:63:THR:CG2	2.63	0.41
45:DO:104:ARG:C	45:DO:106:LEU:H	2.21	0.41
45:DO:2:ILE:O	45:DO:33:ALA:N	2.53	0.41
45:DO:49:ARG:HD3	45:DO:49:ARG:N	2.35	0.41
46:DP:48:PRO:O	46:DP:50:ARG:N	2.52	0.41
48:DR:9:LYS:O	48:DR:10:LEU:CD2	2.68	0.41
49:DS:23:ARG:O	49:DS:24:LEU:C	2.59	0.41
49:DS:66:ALA:O	49:DS:68:GLN:N	2.50	0.41
51:DU:62:ILE:O	51:DU:65:ILE:HB	2.21	0.41
52:DV:39:LEU:O	52:DV:50:PRO:HA	2.20	0.41
35:DA:1601:G:OP2	54:DX:58:HIS:CD2	2.73	0.41
36:DB:105:A:O2'	56:DZ:30:ASN:C	2.58	0.41
56:DZ:44:PHE:CD1	56:DZ:48:PHE:HB2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:5:LEU:HD21	56:DZ:43:GLU:CB	2.51	0.41
56:DZ:77:ASP:O	56:DZ:77:ASP:OD1	2.38	0.41
1:AA:1079:G:C2	1:AA:1080:A:C5	3.08	0.41
1:AA:1128:C:H2'	1:AA:1130:A:C8	2.55	0.41
1:AA:11:G:H2'	1:AA:12:U:H6	1.85	0.41
1:AA:1203:C:OP1	14:AN:3:ARG:CD	2.65	0.41
1:AA:1231:G:H2'	1:AA:1232:U:O4'	2.20	0.41
1:AA:1284:C:H3'	1:AA:1285:A:H8	1.85	0.41
1:AA:1287:A:C5	1:AA:1288:A:N6	2.88	0.41
1:AA:1442:G:C8	1:AA:1442(B):A:N1	2.89	0.41
1:AA:673:G:C6	1:AA:734:G:C6	3.07	0.41
1:AA:69:G:C2	1:AA:70:G:C5	3.08	0.41
1:AA:874:G:O2'	1:AA:875:C:H5'	2.20	0.41
2:AB:100:GLY:CA	2:AB:104:ASN:H	2.34	0.41
2:AB:102:LEU:HD22	2:AB:176:GLU:HB3	2.01	0.41
2:AB:16:HIS:HB3	2:AB:210:SER:OG	2.21	0.41
4:AD:152:SER:O	4:AD:154:ASN:N	2.53	0.41
4:AD:9:CYS:SG	4:AD:22:LYS:CE	3.08	0.41
5:AE:20:GLN:O	5:AE:21:ALA:C	2.58	0.41
1:AA:876:G:H1'	8:AH:11:THR:HG21	2.01	0.41
8:AH:120:THR:C	8:AH:122:ARG:H	2.24	0.41
8:AH:54:ASP:C	8:AH:56:LYS:N	2.74	0.41
8:AH:5:PRO:HA	8:AH:8:ASP:HB3	2.02	0.41
1:AA:1349:A:OP1	9:AI:118:LYS:HB2	2.21	0.41
9:AI:50:LEU:O	9:AI:53:VAL:CG2	2.63	0.41
10:AJ:48:THR:HG23	10:AJ:62:HIS:CA	2.50	0.41
11:AK:115:PRO:C	11:AK:116:HIS:ND1	2.74	0.41
11:AK:26:ASN:O	11:AK:27:ASN:CB	2.67	0.41
11:AK:67:ASP:HA	11:AK:70:LYS:HB3	2.02	0.41
11:AK:84:VAL:HG11	11:AK:95:ILE:HD11	2.00	0.41
12:AL:113:ARG:O	12:AL:122:THR:HG21	2.20	0.41
1:AA:538:G:P	12:AL:115:LYS:HB2	2.60	0.41
12:AL:75:HIS:CD2	12:AL:77:LEU:HG	2.42	0.41
13:AM:83:ASP:OD2	13:AM:84:ILE:N	2.53	0.41
14:AN:33:VAL:HG12	14:AN:34:TYR:H	1.85	0.41
18:AR:26:LEU:HD21	18:AR:42:ARG:NH1	2.36	0.41
26:B0:32:ARG:HB3	26:B0:32:ARG:HE	1.63	0.41
27:B1:41:ARG:HG3	27:B1:41:ARG:NH1	2.30	0.41
31:B5:48:GLU:OE1	31:B5:49:CYS:SG	2.73	0.41
33:B7:29:LYS:NZ	33:B7:32:LYS:NZ	2.68	0.41
33:B7:31:LEU:HD23	33:B7:31:LEU:HA	1.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1153:C:N4	35:BA:1154:G:N1	2.68	0.41
35:BA:1360:A:C6	35:BA:1372:U:C4	3.09	0.41
35:BA:1515:G:N2	35:BA:1516:C:H1'	2.35	0.41
35:BA:1568:G:P	38:BD:63:ARG:HH22	2.44	0.41
35:BA:179:G:C4	35:BA:180:G:C8	3.08	0.41
35:BA:1812:A:C1'	38:BD:46:GLN:HE22	2.33	0.41
35:BA:2068:U:H3	35:BA:2430:A:H2	1.62	0.41
35:BA:2085:C:O2'	35:BA:2086:U:H5'	2.20	0.41
35:BA:210:C:H2'	35:BA:211:A:H8	1.85	0.41
35:BA:2123:G:O2'	35:BA:2124:G:H5'	2.20	0.41
35:BA:2408:U:C2	35:BA:2409:G:C8	3.08	0.41
35:BA:2251:G:C2	35:BA:2450:A:H1'	2.56	0.41
35:BA:2603:G:H4'	35:BA:2603:G:OP2	2.20	0.41
35:BA:2748:A:C2	35:BA:2757:A:C4	3.08	0.41
35:BA:2773:C:OP1	39:BE:164:ARG:HG2	2.20	0.41
35:BA:389:G:C8	35:BA:2413:G:H4'	2.55	0.41
35:BA:754:C:H2'	35:BA:755:C:C6	2.55	0.41
37:BC:89:ALA:HB1	37:BC:152:ILE:O	2.20	0.41
37:BC:74:VAL:HB	37:BC:91:ALA:HB2	2.02	0.41
40:BF:46:ARG:HG3	40:BF:48:THR:CG2	2.50	0.41
40:BF:78:ILE:HG13	40:BF:78:ILE:H	1.65	0.41
41:BG:165:THR:O	41:BG:166:ASP:C	2.58	0.41
41:BG:43:LEU:HD12	41:BG:153:ARG:CD	2.50	0.41
42:BH:58:GLU:O	42:BH:61:HIS:HB2	2.20	0.41
42:BH:87:LEU:HA	42:BH:164:TYR:O	2.20	0.41
43:BI:98:ALA:HB1	43:BI:109:ILE:HD13	2.01	0.41
43:BI:51:ILE:CG2	43:BI:52:ARG:N	2.80	0.41
44:BN:30:ILE:CD1	44:BN:54:VAL:HG21	2.46	0.41
45:BO:87:ILE:CG2	45:BO:91:LEU:HD13	2.51	0.41
46:BP:126:VAL:CA	46:BP:145:PRO:HG2	2.46	0.41
46:BP:33:ARG:O	46:BP:35:HIS:N	2.52	0.41
46:BP:47:ASP:CG	46:BP:49:ARG:HB3	2.40	0.41
47:BQ:24:GLY:CA	47:BQ:101:ARG:HA	2.51	0.41
47:BQ:27:VAL:HG23	56:BZ:81:ARG:HH22	1.85	0.41
35:BA:2470:G:OP1	47:BQ:56:ARG:NH2	2.53	0.41
47:BQ:6:ARG:O	47:BQ:6:ARG:HG3	2.19	0.41
48:BR:9:LYS:HG3	48:BR:43:GLU:OE2	2.20	0.41
49:BS:24:LEU:N	49:BS:24:LEU:HD22	2.36	0.41
49:BS:66:ALA:O	49:BS:68:GLN:N	2.51	0.41
49:BS:87:PHE:CE2	49:BS:88:ASP:O	2.73	0.41
50:BT:48:ILE:HD12	50:BT:48:ILE:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:65:LYS:HG3	50:BT:66:VAL:N	2.35	0.41
50:BT:64:ARG:CB	50:BT:73:GLU:HB3	2.48	0.41
50:BT:88:ILE:CD1	50:BT:88:ILE:N	2.83	0.41
53:BW:10:VAL:O	53:BW:11:ARG:HB2	2.20	0.41
54:BX:40:LYS:HD2	54:BX:41:ASN:CA	2.49	0.41
54:BX:59:VAL:HB	54:BX:60:ARG:H	1.30	0.41
55:BY:14:LEU:HD12	55:BY:23:ARG:O	2.20	0.41
55:BY:28:LYS:O	55:BY:37:VAL:O	2.38	0.41
56:BZ:116:VAL:HG12	56:BZ:117:LEU:CD2	2.48	0.41
56:BZ:28:MET:SD	56:BZ:37:VAL:HG21	2.60	0.41
56:BZ:56:VAL:HG13	56:BZ:57:ILE:N	2.35	0.41
47:BQ:141:GLN:CG	56:BZ:72:ARG:HE	2.32	0.41
1:CA:1003:G:H8	1:CA:1003:G:O5'	2.04	0.41
1:CA:100:C:H2'	1:CA:101:A:C8	2.56	0.41
1:CA:1063:C:H3'	1:CA:1064:G:H2'	2.03	0.41
1:CA:107:G:C2'	1:CA:108:G:H5'	2.50	0.41
1:CA:1113:C:O5'	1:CA:1113:C:H6	2.03	0.41
1:CA:1226:C:C2	13:CM:104:ARG:HA	2.55	0.41
1:CA:1229:A:O2'	1:CA:1230:C:H5'	2.20	0.41
1:CA:1413:A:C2	1:CA:1414:U:C2	3.09	0.41
1:CA:325:A:N6	1:CA:326:G:N1	2.68	0.41
1:CA:477:A:HO2'	1:CA:479:C:H5'	1.82	0.41
1:CA:660:G:C2	1:CA:661:G:C4	3.08	0.41
1:CA:992:U:O2'	1:CA:993:G:P	2.78	0.41
2:CB:164:VAL:HG12	2:CB:165:VAL:N	2.35	0.41
2:CB:11:LEU:O	2:CB:16:His:CE1	2.73	0.41
2:CB:222:ILE:CG2	2:CB:223:ILE:N	2.81	0.41
2:CB:229:VAL:HG12	2:CB:229:VAL:O	2.20	0.41
2:CB:44:LEU:H	2:CB:44:LEU:HG	1.64	0.41
3:CC:14:ILE:HG12	3:CC:15:THR:N	2.35	0.41
3:CC:88:ARG:N	3:CC:101:LEU:HD12	2.35	0.41
8:CH:86:ILE:HG22	8:CH:87:SER:N	2.35	0.41
9:CI:79:LEU:HD13	9:CI:79:LEU:C	2.41	0.41
10:CJ:50:ILE:CD1	10:CJ:50:ILE:N	2.69	0.41
10:CJ:78:ASN:O	10:CJ:80:LYS:N	2.53	0.41
11:CK:86:GLY:H	11:CK:112:THR:CB	2.33	0.41
12:CL:101:VAL:O	12:CL:103:GLY:N	2.53	0.41
12:CL:120:TYR:N	12:CL:120:TYR:CD1	2.88	0.41
12:CL:27:LEU:O	12:CL:28:LYS:C	2.58	0.41
13:CM:68:GLY:O	13:CM:70:LEU:N	2.53	0.41
17:CQ:11:VAL:HG23	17:CQ:20:THR:HB	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CR:65:ILE:HD12	18:CR:66:LEU:H	1.77	0.41
18:CR:79:LEU:HA	18:CR:80:PRO:HD3	1.77	0.41
23:CW:70:C:H2'	23:CW:71:G:C8	2.54	0.41
25:CY:142:LYS:HA	25:CY:145:LYS:HE2	2.02	0.41
25:CY:21:LEU:HD11	25:CY:121:TYR:C	2.40	0.41
25:CY:41:LEU:N	25:CY:41:LEU:HD12	2.22	0.41
31:D5:31:VAL:O	31:D5:32:PRO:C	2.59	0.41
34:D8:56:GLU:C	34:D8:58:ILE:H	2.22	0.41
35:DA:1231:G:H2'	35:DA:1232:G:C8	2.55	0.41
35:DA:1353:A:H4'	38:DD:38:LYS:HZ3	1.84	0.41
35:DA:1416:G:HO2'	35:DA:1417:C:H6	1.60	0.41
35:DA:1433:U:O2	35:DA:1561:G:C2	2.73	0.41
35:DA:1484:G:C3'	35:DA:1485:G:C5'	2.94	0.41
35:DA:1545:A:N7	35:DA:1546:C:C2	2.88	0.41
35:DA:1721:G:C6	35:DA:1739:U:H5'	2.54	0.41
35:DA:1794:U:H2'	35:DA:1795:C:C6	2.55	0.41
35:DA:2320:A:N3	35:DA:2320:A:H2'	2.35	0.41
32:D6:24:GLU:OE2	35:DA:2346:A:O2'	2.39	0.41
35:DA:313:C:H2'	35:DA:314:A:C8	2.50	0.41
35:DA:438:G:H2'	35:DA:440:G:C8	2.55	0.41
35:DA:687:C:H42	35:DA:787:U:H4'	1.85	0.41
35:DA:685:A:N1	35:DA:787:U:H1'	2.35	0.41
35:DA:85:G:N2	35:DA:86:C:H1'	2.35	0.41
35:DA:906:G:H2'	35:DA:907:U:O4'	2.20	0.41
36:DB:24:G:H4'	36:DB:25:A:H8	1.80	0.41
36:DB:71:C:H2'	36:DB:72:G:O4'	2.20	0.41
37:DC:74:VAL:HB	37:DC:91:ALA:HB2	2.03	0.41
38:DD:105:ILE:O	38:DD:105:ILE:HG23	2.20	0.41
38:DD:185:VAL:HG12	38:DD:189:CYS:SG	2.60	0.41
38:DD:53:PHE:CD1	38:DD:219:PRO:O	2.73	0.41
38:DD:29:PRO:O	38:DD:30:GLU:C	2.59	0.41
38:DD:34:VAL:HG22	38:DD:35:LYS:HZ2	1.84	0.41
38:DD:92:ILE:C	38:DD:107:ALA:HB2	2.41	0.41
39:DE:201:THR:CG2	39:DE:203:LYS:HB3	2.50	0.41
39:DE:35:GLN:NE2	39:DE:37:ARG:NH2	2.67	0.41
40:DF:133:ASN:O	40:DF:134:GLY:C	2.58	0.41
40:DF:3:GLU:O	40:DF:19:GLU:CB	2.69	0.41
40:DF:24:LEU:N	40:DF:24:LEU:HD22	2.35	0.41
41:DG:43:LEU:HD11	41:DG:90:LEU:H	1.84	0.41
41:DG:67:LYS:CD	41:DG:67:LYS:H	2.29	0.41
41:DG:96:ARG:C	41:DG:99:MET:HB3	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DN:44:PRO:C	44:DN:46:VAL:N	2.72	0.41
45:DO:8:LEU:N	45:DO:8:LEU:HD22	2.36	0.41
46:DP:59:LEU:HA	46:DP:61:ARG:HD2	2.02	0.41
47:DQ:115:MET:CE	47:DQ:133:ARG:HH21	2.33	0.41
49:DS:24:LEU:HD22	49:DS:24:LEU:N	2.34	0.41
49:DS:69:VAL:CG1	49:DS:70:GLY:N	2.83	0.41
49:DS:77:ALA:O	49:DS:79:ALA:N	2.53	0.41
50:DT:35:LYS:CE	50:DT:41:ARG:HG3	2.50	0.41
51:DU:103:PRO:O	51:DU:104:GLN:C	2.59	0.41
55:DY:8:LYS:HZ1	55:DY:74:PRO:HD3	1.86	0.41
56:DZ:108:PRO:HB2	56:DZ:144:LEU:H	1.83	0.41
56:DZ:151:HIS:CA	56:DZ:170:THR:HA	2.47	0.41
56:DZ:22:GLY:C	56:DZ:41:LEU:HD21	2.41	0.41
1:AA:1007:C:H2'	1:AA:1008:C:C5	2.55	0.41
1:AA:1023:G:C2'	1:AA:1024:G:H5'	2.50	0.41
1:AA:1210:C:H4'	1:AA:1214:C:N4	2.35	0.41
1:AA:1227:A:N7	1:AA:1228:C:C2	2.88	0.41
1:AA:1248:A:H2'	1:AA:1249:C:C5'	2.49	0.41
1:AA:55:A:N7	1:AA:56:U:C4	2.88	0.41
1:AA:666:G:O2'	1:AA:667:G:H5'	2.20	0.41
1:AA:690:G:C6	1:AA:691:G:C2	3.09	0.41
1:AA:778:G:C2'	1:AA:779:C:H5'	2.50	0.41
1:AA:926:G:C6	1:AA:1505:G:C6	3.07	0.41
1:AA:935:A:O2'	1:AA:936:C:H5'	2.21	0.41
2:AB:132:LYS:HA	2:AB:135:GLN:NE2	2.36	0.41
2:AB:229:VAL:O	2:AB:229:VAL:HG12	2.20	0.41
3:AC:11:ARG:O	3:AC:13:GLY:N	2.53	0.41
3:AC:95:THR:C	3:AC:97:LYS:N	2.74	0.41
4:AD:154:ASN:HB2	4:AD:159:ARG:HH21	1.84	0.41
5:AE:101:ILE:HG12	5:AE:118:ILE:O	2.19	0.41
5:AE:36:ASP:CG	5:AE:37:ARG:N	2.73	0.41
5:AE:68:GLU:OE2	5:AE:70:PRO:HD3	2.21	0.41
7:AG:40:ALA:O	7:AG:41:ARG:C	2.59	0.41
7:AG:41:ARG:CG	7:AG:41:ARG:HH11	2.32	0.41
8:AH:54:ASP:C	8:AH:56:LYS:H	2.23	0.41
8:AH:4:ASP:OD2	8:AH:7:ALA:CB	2.68	0.41
10:AJ:64:GLU:HG3	10:AJ:64:GLU:O	2.19	0.41
12:AL:60:LEU:CD2	12:AL:64:TYR:O	2.67	0.41
14:AN:53:LEU:HA	14:AN:53:LEU:HD23	1.88	0.41
17:AQ:11:VAL:HG23	17:AQ:20:THR:CG2	2.51	0.41
17:AQ:29:HIS:CE1	17:AQ:31:LEU:H	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:AQ:60:ILE:HB	17:AQ:74:LEU:HD23	2.02	0.41
17:AQ:68:ARG:H	17:AQ:70:ARG:HH12	1.68	0.41
17:AQ:58:GLU:O	17:AQ:74:LEU:CB	2.67	0.41
1:AA:1319:A:OP2	19:AS:5:LEU:HD23	2.20	0.41
1:AA:1223:C:P	19:AS:78:ARG:HH12	2.43	0.41
25:AY:6:LEU:C	25:AY:8:ALA:H	2.23	0.41
26:B0:38:VAL:HG12	26:B0:39:ARG:N	2.36	0.41
31:B5:16:ARG:HG2	31:B5:16:ARG:NH1	2.26	0.41
32:B6:10:LEU:O	32:B6:11:LEU:HD22	2.21	0.41
34:B8:49:VAL:O	34:B8:50:LEU:HB3	2.20	0.41
35:BA:942:G:H1'	35:BA:1189:A:C2	2.55	0.41
35:BA:1417:C:H2'	35:BA:1418:G:H5'	2.01	0.41
35:BA:1493:C:O2	35:BA:1493:C:H2'	2.21	0.41
35:BA:1767:C:O2'	35:BA:1768:U:H5'	2.20	0.41
35:BA:1839:G:C4	35:BA:1840:G:C8	3.08	0.41
35:BA:1891:G:C2	35:BA:1892:C:C2	3.08	0.41
35:BA:1665:A:N1	35:BA:1996:C:N4	2.69	0.41
31:B5:19:ARG:HA	35:BA:2046:G:O5'	2.20	0.41
35:BA:2078:C:C2	35:BA:2079:U:C6	3.08	0.41
35:BA:227:A:C2	35:BA:2407:G:C1'	2.99	0.41
34:B8:31:HIS:CG	35:BA:2419:U:O4	2.74	0.41
35:BA:2464:C:N4	35:BA:2487:G:N1	2.68	0.41
35:BA:2753:A:H2	35:BA:2754:U:N3	2.18	0.41
35:BA:2787:C:O2	39:BE:61:ARG:NH1	2.53	0.41
35:BA:2832:U:C2	35:BA:2834:G:N2	2.88	0.41
35:BA:28:A:C8	35:BA:513:A:C5	3.08	0.41
35:BA:292:C:O2	35:BA:292:C:C2'	2.69	0.41
35:BA:374:A:H2'	35:BA:375:C:H5'	2.02	0.41
27:B1:19:GLN:HB3	35:BA:380:U:O2'	2.21	0.41
35:BA:520:G:O2'	35:BA:521:G:H5'	2.19	0.41
35:BA:604:G:O6	35:BA:625:G:C6	2.72	0.41
35:BA:742:G:N1	35:BA:756:C:N4	2.68	0.41
35:BA:811:U:O2	35:BA:1251:C:C6	2.73	0.41
36:BB:32:C:C2	36:BB:51:G:N2	2.89	0.41
38:BD:77:ALA:CB	38:BD:97:TYR:CD1	3.03	0.41
38:BD:9:TYR:C	38:BD:10:THR:CG2	2.88	0.41
35:BA:2575:C:C5'	39:BE:144:ARG:HD3	2.46	0.41
39:BE:48:GLN:O	39:BE:49:LEU:HD13	2.20	0.41
39:BE:49:LEU:O	39:BE:78:LEU:CA	2.69	0.41
39:BE:61:ARG:CG	39:BE:62:PRO:N	2.84	0.41
42:BH:158:HIS:HE1	42:BH:168:PRO:HB2	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BH:54:ARG:HB2	42:BH:61:HIS:HD2	1.85	0.41
43:BI:7:GLU:O	43:BI:9:LEU:HG	2.20	0.41
44:BN:95:PRO:O	44:BN:96:GLU:C	2.59	0.41
46:BP:140:ALA:O	46:BP:141:ALA:CB	2.67	0.41
49:BS:13:ARG:N	49:BS:13:ARG:CD	2.70	0.41
50:BT:28:VAL:HG11	50:BT:46:GLU:CD	2.41	0.41
51:BU:65:ILE:HG22	51:BU:66:ASN:N	2.35	0.41
52:BV:22:VAL:HG21	52:BV:96:ILE:CG1	2.51	0.41
53:BW:62:HIS:O	53:BW:64:MET:HG3	2.20	0.41
55:BY:81:LYS:HA	55:BY:82:PRO:HD3	1.84	0.41
47:BQ:141:GLN:OE1	56:BZ:70:LEU:HB3	2.20	0.41
1:CA:1011:G:H2'	1:CA:1012:U:C5'	2.50	0.41
1:CA:1128:C:H2'	1:CA:1130:A:C8	2.55	0.41
1:CA:1346:A:N6	1:CA:1374:A:H3'	2.35	0.41
1:CA:1422:G:C2	1:CA:1423:G:C8	3.08	0.41
1:CA:1477:C:O2'	1:CA:1478:C:H5'	2.20	0.41
1:CA:781:A:O2'	1:CA:1522:U:O2	2.36	0.41
1:CA:315:A:H5''	1:CA:317:G:OP2	2.21	0.41
1:CA:321:A:H4'	1:CA:1436:U:C4'	2.50	0.41
1:CA:450:G:OP1	1:CA:452:A:OP2	2.39	0.41
1:CA:607:A:C4	1:CA:608:A:C8	3.09	0.41
1:CA:729:A:H2'	1:CA:730:G:H8	1.85	0.41
1:CA:858:G:O6	1:CA:869:G:C8	2.73	0.41
2:CB:166:ASP:O	2:CB:167:PRO:O	2.38	0.41
3:CC:172:ARG:NH1	3:CC:172:ARG:HB3	2.36	0.41
3:CC:22:TRP:CZ3	3:CC:32:LEU:HB3	2.55	0.41
4:CD:101:LEU:O	4:CD:104:VAL:HB	2.20	0.41
4:CD:24:GLU:O	4:CD:27:TYR:N	2.50	0.41
4:CD:18:LYS:HZ1	4:CD:31:CYS:HB3	1.83	0.41
5:CE:131:ILE:O	5:CE:132:ALA:C	2.59	0.41
5:CE:72:GLN:O	5:CE:73:ASN:CB	2.69	0.41
5:CE:87:SER:OG	5:CE:125:SER:CB	2.67	0.41
8:CH:48:TYR:O	8:CH:49:GLU:HB3	2.20	0.41
9:CI:116:LYS:O	9:CI:119:ALA:N	2.54	0.41
9:CI:126:SER:O	9:CI:128:ARG:HD3	2.21	0.41
10:CJ:4:ILE:N	10:CJ:4:ILE:HD12	2.36	0.41
10:CJ:8:LEU:H	10:CJ:8:LEU:HD12	1.85	0.41
14:CN:27:CYS:CB	14:CN:43:CYS:SG	3.02	0.41
17:CQ:58:GLU:O	17:CQ:74:LEU:CB	2.69	0.41
1:CA:1226:C:H5'	19:CS:80:TYR:CE2	2.56	0.41
20:CT:78:ALA:O	20:CT:79:ARG:C	2.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:CY:25:LEU:C	25:CY:27:GLY:H	2.23	0.41
28:D2:32:LEU:HD22	28:D2:44:LEU:HD21	2.01	0.41
31:D5:31:VAL:O	31:D5:32:PRO:O	2.37	0.41
33:D7:31:LEU:CD2	33:D7:42:LEU:HD22	2.50	0.41
34:D8:26:LYS:HZ3	34:D8:47:LYS:HD3	1.85	0.41
35:DA:1106:A:H2'	35:DA:1107:G:C8	2.55	0.41
35:DA:1681:G:H1'	35:DA:1763:G:H5'	2.03	0.41
35:DA:1947:C:H42	35:DA:1959:G:H1	1.67	0.41
35:DA:195:A:O5'	35:DA:196:A:H4'	2.20	0.41
35:DA:2014:A:H2'	35:DA:2015:A:C8	2.55	0.41
35:DA:2288:A:H2	35:DA:2325:G:C8	2.38	0.41
35:DA:2419:U:H2'	35:DA:2420:C:C6	2.56	0.41
35:DA:2545:G:C2	35:DA:2546:U:C2	3.08	0.41
35:DA:270:A:OP2	35:DA:271(X):G:N2	2.49	0.41
35:DA:2820:A:C8	39:DE:191:PRO:CB	2.94	0.41
35:DA:331:A:O2'	35:DA:332:A:OP1	2.36	0.41
35:DA:542:C:N3	35:DA:543:C:N4	2.55	0.41
35:DA:548:A:N3	35:DA:548:A:O2'	2.52	0.41
34:D8:18:ALA:HB2	35:DA:628:G:H5''	2.03	0.41
35:DA:769:G:H5'	35:DA:1379:A:H61	1.84	0.41
35:DA:687:C:N3	35:DA:788:A:H5'	2.35	0.41
35:DA:863:A:H2	35:DA:914:C:N4	2.18	0.41
37:DC:51:PRO:HB3	37:DC:204:ALA:CB	2.50	0.41
38:DD:13:ARG:HD2	38:DD:13:ARG:HA	1.95	0.41
38:DD:142:VAL:CG2	38:DD:192:THR:C	2.86	0.41
38:DD:48:ARG:NH1	38:DD:48:ARG:CG	2.81	0.41
39:DE:61:ARG:HG2	39:DE:62:PRO:CD	2.46	0.41
41:DG:63:ILE:HB	41:DG:141:PHE:CE1	2.56	0.41
41:DG:48:GLU:HG2	41:DG:49:ASP:N	2.29	0.41
41:DG:53:LEU:HD22	41:DG:56:ALA:HB2	2.02	0.41
41:DG:90:LEU:HD12	41:DG:90:LEU:N	2.35	0.41
43:DI:62:LYS:HD3	43:DI:62:LYS:C	2.40	0.41
44:DN:27:ALA:O	44:DN:30:ILE:N	2.53	0.41
48:DR:21:TYR:HB3	48:DR:47:PHE:CD2	2.55	0.41
48:DR:28:LEU:HD12	48:DR:29:LEU:HD13	2.02	0.41
50:DT:100:TYR:HD1	50:DT:100:TYR:H	1.68	0.41
53:DW:62:HIS:O	53:DW:63:ASP:C	2.59	0.41
53:DW:96:ILE:HG13	53:DW:97:LYS:N	2.35	0.41
55:DY:20:TYR:CZ	55:DY:42:VAL:HA	2.55	0.41
55:DY:28:LYS:CB	55:DY:39:VAL:H	2.32	0.41
55:DY:47:LYS:HD2	55:DY:47:LYS:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:DY:62:GLU:OE1	55:DY:62:GLU:HA	2.20	0.41
1:AA:10:A:H2'	1:AA:11:G:C8	2.52	0.41
1:AA:110:C:C4	1:AA:111:G:C5	3.09	0.41
1:AA:1365:G:H2'	1:AA:1366:C:C5'	2.50	0.41
1:AA:1064:G:OP2	1:AA:1386:G:H4'	2.20	0.41
1:AA:1511:G:C5	1:AA:1512:U:C5	3.09	0.41
1:AA:516:U:C4	1:AA:517:G:C6	3.08	0.41
1:AA:770:C:HO2'	1:AA:771:G:H5'	1.83	0.41
1:AA:939:G:H2'	1:AA:940:C:C6	2.55	0.41
2:AB:95:GLN:HG3	2:AB:147:LYS:O	2.20	0.41
2:AB:204:ASN:HB3	2:AB:210:SER:HB3	2.01	0.41
2:AB:87:ARG:O	2:AB:223:ILE:HD11	2.21	0.41
4:AD:52:SER:O	4:AD:55:ALA:N	2.54	0.41
4:AD:63:LYS:O	4:AD:64:LEU:C	2.59	0.41
5:AE:11:ILE:HD12	5:AE:31:LEU:CD2	2.48	0.41
5:AE:126:ARG:NH1	5:AE:126:ARG:HG3	2.35	0.41
8:AH:102:ARG:N	8:AH:102:ARG:HE	2.19	0.41
8:AH:107:LEU:HD23	8:AH:107:LEU:O	2.20	0.41
7:AG:16:LEU:CD1	9:AI:42:ARG:HA	2.51	0.41
9:AI:82:ALA:HA	9:AI:85:LEU:HD21	2.02	0.41
10:AJ:4:ILE:HD12	10:AJ:4:ILE:N	2.35	0.41
11:AK:106:LYS:HB3	11:AK:107:SER:H	1.57	0.41
11:AK:122:LYS:O	11:AK:126:ARG:N	2.53	0.41
11:AK:30:VAL:O	11:AK:30:VAL:HG23	2.20	0.41
1:AA:523:A:N6	12:AL:53:ARG:NH2	2.69	0.41
13:AM:16:ASP:N	13:AM:16:ASP:OD2	2.53	0.41
15:AO:66:LEU:O	15:AO:67:LEU:C	2.59	0.41
17:AQ:3:LYS:O	17:AQ:4:LYS:C	2.59	0.41
19:AS:6:LYS:HG2	19:AS:7:LYS:HD2	2.02	0.41
20:AT:82:SER:O	20:AT:86:ARG:HB2	2.19	0.41
25:AY:92:PRO:N	25:AY:101:ILE:HG23	2.35	0.41
25:AY:148:HIS:C	25:AY:149:LEU:HD23	2.41	0.41
27:B1:66:HIS:N	27:B1:66:HIS:CD2	2.85	0.41
27:B1:90:ILE:C	27:B1:93:GLU:OE2	2.59	0.41
32:B6:47:THR:CG2	32:B6:48:VAL:N	2.84	0.41
34:B8:56:GLU:CD	34:B8:56:GLU:N	2.74	0.41
35:BA:109:G:C4	35:BA:110:G:C8	3.08	0.41
35:BA:1112:G:O2'	35:BA:1113:U:H6	2.03	0.41
35:BA:818:G:C2	35:BA:1190:G:C6	3.08	0.41
35:BA:1297:C:O2'	35:BA:1302:A:N1	2.54	0.41
35:BA:1342:A:C2	35:BA:1345:C:C5	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1360:A:C2	35:BA:1372:U:C2	3.08	0.41
35:BA:1778:U:C5	35:BA:1784:A:N3	2.88	0.41
35:BA:1923:U:H2'	35:BA:1924:C:C6	2.56	0.41
35:BA:1947:C:H42	35:BA:1959:G:H1	1.68	0.41
35:BA:2023:G:H2'	35:BA:2024:G:H8	1.84	0.41
35:BA:2633:G:C2	35:BA:2634:G:C4	3.09	0.41
35:BA:2812:G:H2'	35:BA:2813:A:H8	1.86	0.41
35:BA:2861:G:C2	35:BA:2862:G:C8	3.09	0.41
35:BA:349:G:H2'	35:BA:349:G:N3	2.35	0.41
35:BA:976:C:H2'	35:BA:977:G:C8	2.56	0.41
37:BC:124:GLY:O	37:BC:125:SER:CB	2.67	0.41
38:BD:223:GLY:HA2	38:BD:226:MET:SD	2.60	0.41
39:BE:117:MET:HE1	39:BE:124:GLY:HA3	2.02	0.41
39:BE:168:MET:O	39:BE:170:LEU:HD12	2.20	0.41
39:BE:4:ILE:HD12	39:BE:5:LEU:H	1.86	0.41
41:BG:115:ARG:HH12	41:BG:136:ARG:HG3	1.85	0.41
41:BG:88:ILE:HG23	41:BG:90:LEU:HD13	2.02	0.41
41:BG:97:ASP:C	41:BG:99:MET:N	2.72	0.41
42:BH:116:GLU:HG2	42:BH:117:PRO:CD	2.48	0.41
42:BH:98:LEU:HD22	42:BH:125:VAL:HB	2.02	0.41
43:BI:83:ALA:CB	43:BI:88:ILE:HG12	2.45	0.41
44:BN:128:HIS:CD2	44:BN:131:GLN:HB2	2.55	0.41
44:BN:46:VAL:CG1	44:BN:47:ALA:N	2.66	0.41
44:BN:51:PHE:O	44:BN:52:VAL:C	2.59	0.41
46:BP:100:LEU:O	46:BP:105:LEU:O	2.38	0.41
46:BP:147:LEU:HB2	46:BP:148:LEU:H	1.58	0.41
50:BT:30:VAL:HA	50:BT:43:GLN:O	2.21	0.41
51:BU:4:ALA:O	51:BU:5:LYS:O	2.39	0.41
52:BV:63:GLY:O	52:BV:64:HIS:CB	2.69	0.41
51:BU:40:PHE:HD1	52:BV:78:LYS:NZ	2.18	0.41
54:BX:53:LYS:CE	54:BX:55:ASN:HD21	2.33	0.41
55:BY:39:VAL:HB	55:BY:40:GLU:H	1.61	0.41
55:BY:86:ARG:CG	55:BY:87:LYS:N	2.82	0.41
56:BZ:122:ARG:CG	56:BZ:123:ASP:OD1	2.68	0.41
56:BZ:166:SER:HB2	56:BZ:167:PRO:C	2.41	0.41
1:CA:125:U:H2'	1:CA:126:G:H8	1.85	0.41
1:CA:1437:C:C2'	1:CA:1438:G:O5'	2.68	0.41
1:CA:180:U:H2'	1:CA:181:G:C5'	2.49	0.41
1:CA:296:U:H2'	1:CA:297:G:O4'	2.21	0.41
1:CA:298:A:H2'	1:CA:299:G:O4'	2.20	0.41
1:CA:327:A:C2	1:CA:329:A:C4	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:585:G:O2'	1:CA:879:C:OP1	2.39	0.41
1:CA:678:U:N3	1:CA:713:G:N2	2.68	0.41
1:CA:707:C:O2	1:CA:707:C:C2'	2.67	0.41
1:CA:837:G:C2	1:CA:850:U:O2	2.74	0.41
1:CA:874:G:C4	1:CA:875:C:C5	3.08	0.41
1:CA:969:A:O2'	1:CA:970:C:H5'	2.20	0.41
1:CA:973:G:H4'	10:CJ:54:PHE:O	2.20	0.41
2:CB:132:LYS:HA	2:CB:135:GLN:NE2	2.35	0.41
2:CB:187:LEU:HD23	2:CB:202:PRO:O	2.20	0.41
4:CD:176:LEU:HD12	4:CD:177:ASP:N	2.34	0.41
5:CE:57:LYS:CE	5:CE:61:TYR:HE2	2.31	0.41
6:CF:17:SER:O	6:CF:21:LEU:CD2	2.68	0.41
6:CF:42:GLU:C	6:CF:44:GLY:H	2.23	0.41
7:CG:64:GLN:C	7:CG:66:VAL:H	2.24	0.41
1:CA:1347:G:N3	9:CI:107:ARG:NH2	2.68	0.41
9:CI:39:GLY:O	9:CI:41:VAL:N	2.54	0.41
10:CJ:50:ILE:HA	10:CJ:60:ARG:CG	2.47	0.41
12:CL:56:ALA:O	12:CL:68:ALA:N	2.53	0.41
16:CP:39:TYR:OH	16:CP:41:PRO:HB3	2.21	0.41
18:CR:74:ARG:CZ	18:CR:81:PHE:HA	2.50	0.41
20:CT:38:LYS:HA	20:CT:41:ILE:HG12	2.02	0.41
21:CU:21:TYR:HD1	21:CU:21:TYR:N	2.19	0.41
25:CY:160:GLU:O	25:CY:163:LYS:N	2.53	0.41
27:D1:23:LYS:CB	27:D1:23:LYS:NZ	2.83	0.41
27:D1:11:ARG:HG3	27:D1:61:ARG:C	2.41	0.41
30:D4:21:VAL:C	30:D4:23:GLU:N	2.73	0.41
31:D5:51:TYR:O	31:D5:56:LYS:HE3	2.21	0.41
35:DA:127:A:H5''	35:DA:128:C:C6	2.54	0.41
35:DA:139:G:H1	35:DA:142(A):C:N4	2.17	0.41
35:DA:1515:G:N2	35:DA:1516:C:H1'	2.35	0.41
35:DA:1550:C:O2'	35:DA:1551:C:H5'	2.21	0.41
35:DA:1782:C:H2'	35:DA:1783:A:H5'	2.03	0.41
35:DA:1832:C:H2'	35:DA:1833:U:O4'	2.21	0.41
35:DA:2127:G:O3'	35:DA:2128:C:C4'	2.68	0.41
35:DA:2287:A:N6	35:DA:2344:U:N3	2.68	0.41
35:DA:2352:A:H2'	35:DA:2353:G:O4'	2.21	0.41
35:DA:2469:A:O2'	47:DQ:56:ARG:CG	2.69	0.41
35:DA:2636:U:H4'	39:DE:80:GLU:OE2	2.21	0.41
35:DA:2730:C:H2'	35:DA:2731:G:C8	2.55	0.41
35:DA:2790:A:N3	35:DA:2790:A:C2'	2.83	0.41
35:DA:359:A:H2'	35:DA:360:G:O4'	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:36:G:O2'	35:DA:37:C:H5'	2.20	0.41
35:DA:587:C:H4'	35:DA:588:U:O5'	2.20	0.41
35:DA:593:G:H2'	35:DA:594:U:H6	1.86	0.41
34:D8:18:ALA:HB3	35:DA:651:G:H4'	2.01	0.41
35:DA:671:C:H2'	35:DA:671:C:O2	2.19	0.41
35:DA:704:G:C2	35:DA:726:G:C4	3.09	0.41
35:DA:727:A:H3'	35:DA:728:G:C8	2.55	0.41
35:DA:845:G:C8	35:DA:845:G:OP2	2.73	0.41
36:DB:102:A:H2'	36:DB:102:A:N3	2.35	0.41
37:DC:44:HIS:CD2	37:DC:175:VAL:HA	2.55	0.41
38:DD:113:VAL:C	38:DD:115:GLN:N	2.73	0.41
38:DD:130:ALA:C	38:DD:131:LEU:HG	2.39	0.41
39:DE:103:ASP:CG	39:DE:168:MET:HG2	2.40	0.41
39:DE:169:ASN:OD1	39:DE:201:THR:CG2	2.67	0.41
39:DE:47:VAL:CG1	39:DE:49:LEU:HD21	2.49	0.41
41:DG:108:ASN:CA	41:DG:112:PRO:HD2	2.50	0.41
41:DG:166:ASP:OD1	41:DG:166:ASP:O	2.39	0.41
41:DG:27:ASN:ND2	41:DG:29:TRP:HD1	2.19	0.41
41:DG:59:GLU:CD	41:DG:60:LEU:HD23	2.41	0.41
42:DH:16:SER:HB2	42:DH:27:LYS:O	2.19	0.41
44:DN:26:LEU:O	44:DN:27:ALA:C	2.58	0.41
44:DN:39:ARG:NH1	44:DN:39:ARG:HG3	2.35	0.41
45:DO:111:PHE:CB	45:DO:114:ILE:HD13	2.37	0.41
46:DP:115:LEU:HB3	46:DP:131:SER:OG	2.19	0.41
47:DQ:10:ARG:HB3	47:DQ:11:LYS:H	1.36	0.41
47:DQ:34:LEU:HD11	47:DQ:129:THR:CG2	2.50	0.41
48:DR:44:LEU:CD1	48:DR:48:VAL:HG23	2.50	0.41
49:DS:58:LEU:HD23	49:DS:65:VAL:HG13	2.02	0.41
50:DT:23:ARG:HH21	50:DT:120:ARG:HD3	1.85	0.41
51:DU:10:ARG:O	51:DU:11:ARG:C	2.58	0.41
51:DU:51:LYS:CA	51:DU:51:LYS:HE2	2.48	0.41
51:DU:65:ILE:HG22	51:DU:66:ASN:N	2.34	0.41
53:DW:10:VAL:HG23	53:DW:101:SER:O	2.21	0.41
53:DW:59:VAL:C	53:DW:61:ASN:H	2.24	0.41
55:DY:76:CYS:HB3	55:DY:96:ILE:CD1	2.50	0.41
55:DY:76:CYS:CB	55:DY:77:PRO:CD	2.98	0.41
1:AA:1277:C:H4'	1:AA:1282:C:O2	2.21	0.41
1:AA:256:U:H2'	1:AA:257:G:H8	1.84	0.41
1:AA:329:A:C5	1:AA:332:G:C6	3.09	0.41
1:AA:336:C:HO2'	1:AA:337:C:H5'	1.84	0.41
1:AA:416:G:C6	1:AA:417:C:C4	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:475:G:H2'	1:AA:476:G:C8	2.55	0.41
1:AA:528:C:H2'	1:AA:529:G:C8	2.50	0.41
1:AA:702:A:C3'	1:AA:703:G:H5'	2.47	0.41
2:AB:158:LEU:O	2:AB:159:PRO:C	2.58	0.41
2:AB:17:PHE:O	2:AB:18:GLY:O	2.38	0.41
2:AB:211:ILE:O	2:AB:212:GLN:C	2.57	0.41
2:AB:44:LEU:HD12	2:AB:45:GLN:N	2.36	0.41
4:AD:60:GLU:HG3	4:AD:198:VAL:HG13	2.03	0.41
7:AG:64:GLN:C	7:AG:66:VAL:H	2.24	0.41
9:AI:95:LYS:CD	9:AI:96:LEU:N	2.77	0.41
10:AJ:8:LEU:H	10:AJ:8:LEU:HD12	1.84	0.41
11:AK:104:GLN:O	11:AK:106:LYS:N	2.53	0.41
11:AK:60:ALA:O	11:AK:62:GLN:N	2.53	0.41
12:AL:85:ILE:HA	12:AL:85:ILE:HD12	1.82	0.41
13:AM:100:GLY:C	13:AM:101:GLN:HG3	2.41	0.41
16:AP:1:MET:O	16:AP:24:ALA:HB2	2.21	0.41
16:AP:76:GLN:CG	16:AP:76:GLN:O	2.68	0.41
17:AQ:67:LYS:HG2	17:AQ:68:ARG:N	2.34	0.41
18:AR:40:LEU:O	18:AR:43:PHE:N	2.53	0.41
18:AR:76:LEU:C	18:AR:78:LEU:H	2.23	0.41
19:AS:36:ARG:NH2	19:AS:72:GLY:O	2.54	0.41
20:AT:46:GLU:O	20:AT:46:GLU:HG2	2.21	0.41
23:AW:57:C:H6	23:AW:57:C:O5'	2.04	0.41
25:AY:112:LYS:CA	25:AY:116:ARG:HG3	2.50	0.41
25:AY:140:LEU:HD21	25:AY:158:GLU:N	2.36	0.41
25:AY:36:ALA:C	25:AY:38:LEU:N	2.74	0.41
26:B0:24:LYS:HG3	35:BA:2355:C:H4'	2.01	0.41
26:B0:50:ASN:ND2	26:B0:81:VAL:O	2.54	0.41
27:B1:94:LEU:HD13	27:B1:94:LEU:C	2.40	0.41
31:B5:40:LYS:HG2	31:B5:46:CYS:HB2	2.01	0.41
35:BA:1037:G:N2	35:BA:1038:C:C2	2.88	0.41
35:BA:1050:A:C2	35:BA:2751:G:C4	3.09	0.41
35:BA:1287:A:OP1	48:BR:105:ARG:O	2.39	0.41
35:BA:1289:C:C2	35:BA:1290:C:C5	3.09	0.41
35:BA:1353:A:C2	35:BA:1354:A:C2	3.09	0.41
35:BA:1365:A:H2'	35:BA:1366:A:H8	1.85	0.41
35:BA:1376:C:O2'	35:BA:1377:G:H5'	2.20	0.41
35:BA:1778:U:C5	35:BA:1784:A:C2	3.09	0.41
35:BA:1814:G:H2'	35:BA:1815:A:C8	2.56	0.41
35:BA:1950:G:O5'	35:BA:1950:G:H8	2.02	0.41
35:BA:2120:G:O2'	35:BA:2121:G:H5'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2290:G:C8	35:BA:2290:G:H5'	2.44	0.41
35:BA:2334:G:N3	49:BS:15:ARG:NH1	2.69	0.41
35:BA:2476:A:H2'	35:BA:2477:C:C5'	2.29	0.41
35:BA:2620:C:C4	35:BA:2621:A:N7	2.89	0.41
35:BA:2692:C:H2'	35:BA:2693:A:C8	2.54	0.41
35:BA:336:C:O2'	35:BA:337:C:H5'	2.21	0.41
35:BA:485:C:C2	35:BA:496:G:N2	2.89	0.41
35:BA:543:C:H6	35:BA:547:A:H8	1.67	0.41
35:BA:632:A:C2	35:BA:2403:C:H1'	2.55	0.41
35:BA:645:C:H2'	35:BA:645:C:O2	2.20	0.41
35:BA:58:G:N3	35:BA:70:G:C2	2.88	0.41
35:BA:729:G:H3'	35:BA:729:G:N3	2.35	0.41
35:BA:858:U:H1'	35:BA:2268:A:H2'	2.03	0.41
35:BA:974:G:C4	35:BA:989:G:C2	3.08	0.41
38:BD:136:ILE:HA	38:BD:137:PRO:HD3	1.94	0.41
38:BD:48:ARG:CG	38:BD:48:ARG:NH1	2.84	0.41
38:BD:52:ARG:CB	38:BD:53:PHE:CE2	3.04	0.41
35:BA:1491:G:H5'	38:BD:99:ASP:OD1	2.21	0.41
39:BE:3:GLY:CA	39:BE:81:ILE:HG21	2.49	0.41
40:BF:25:PRO:HB3	40:BF:119:ARG:CG	2.50	0.41
43:BI:125:GLU:OE1	43:BI:125:GLU:HA	2.20	0.41
44:BN:128:HIS:HD2	44:BN:131:GLN:HB2	1.85	0.41
48:BR:116:LEU:CD2	48:BR:117:VAL:HG12	2.51	0.41
48:BR:29:LEU:HG	48:BR:79:LEU:HD23	2.01	0.41
49:BS:73:LEU:O	49:BS:74:ALA:C	2.59	0.41
49:BS:82:ILE:CG2	49:BS:83:LYS:N	2.84	0.41
54:BX:18:TYR:HA	54:BX:21:PHE:CG	2.54	0.41
54:BX:65:ARG:O	54:BX:66:LEU:HB2	2.20	0.41
54:BX:82:GLN:C	54:BX:85:PRO:HD2	2.41	0.41
1:CA:1146:A:C2'	1:CA:1147:C:O5'	2.69	0.41
1:CA:1226:C:H2'	13:CM:103:THR:CB	2.51	0.41
1:CA:1442:G:C5	1:CA:1442(B):A:C2	3.07	0.41
1:CA:1445:C:H2'	1:CA:1446:U:C6	2.42	0.41
1:CA:149:A:O2'	1:CA:150:C:C6	2.73	0.41
1:CA:328:C:HO2'	1:CA:329:A:P	2.42	0.41
1:CA:437:U:H2'	1:CA:438:G:H5'	2.03	0.41
1:CA:502:G:C6	1:CA:503:C:C4	3.08	0.41
1:CA:541:G:H2'	1:CA:542:G:C8	2.56	0.41
1:CA:570:G:C6	1:CA:873:A:C2	3.08	0.41
1:CA:739:C:O2'	1:CA:740:U:H5'	2.21	0.41
1:CA:777:A:C2	1:CA:778:G:C4	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:662:G:O2'	1:CA:836:G:H5''	2.20	0.41
1:CA:909:A:H2'	1:CA:910:C:O4'	2.19	0.41
2:CB:188:ALA:HB1	2:CB:192:SER:HB2	2.02	0.41
2:CB:239:VAL:O	2:CB:240:GLN:HB3	2.20	0.41
1:CA:1100:C:OP2	2:CB:96:ARG:HG2	2.21	0.41
3:CC:134:ILE:HG23	3:CC:151:VAL:HG11	2.03	0.41
3:CC:121:ALA:HB1	3:CC:188:LEU:O	2.19	0.41
1:CA:1190:G:OP1	3:CC:5:ILE:HG23	2.18	0.41
4:CD:131:ARG:O	4:CD:132:ARG:O	2.39	0.41
4:CD:202:LEU:O	4:CD:203:VAL:C	2.56	0.41
4:CD:50:ARG:HD2	4:CD:50:ARG:C	2.41	0.41
5:CE:102:ALA:CB	5:CE:106:PRO:HG2	2.50	0.41
6:CF:14:LEU:HD22	6:CF:15:ASP:N	2.35	0.41
7:CG:111:ARG:HB3	7:CG:112:PRO:HD2	2.03	0.41
7:CG:155:ARG:HG3	7:CG:155:ARG:HH11	1.86	0.41
10:CJ:51:ARG:HG3	10:CJ:60:ARG:C	2.41	0.41
11:CK:72:ALA:O	11:CK:77:MET:HB2	2.20	0.41
13:CM:108:ARG:O	13:CM:109:THR:C	2.58	0.41
17:CQ:27:PHE:C	17:CQ:27:PHE:CD1	2.94	0.41
17:CQ:29:HIS:CG	17:CQ:30:PRO:HD2	2.56	0.41
12:CL:10:LEU:HB3	17:CQ:32:TYR:CE1	2.55	0.41
17:CQ:43:LEU:HD23	17:CQ:43:LEU:HA	1.85	0.41
17:CQ:76:LEU:HD12	17:CQ:77:VAL:N	2.33	0.41
22:CV:40:C:O2	22:CV:41:C:C6	2.74	0.41
25:CY:127:VAL:HA	25:CY:130:ARG:HG3	2.01	0.41
25:CY:62:ASP:HB2	25:CY:63:PRO:HD2	2.02	0.41
26:D0:69:PHE:CD2	26:D0:79:VAL:HG22	2.56	0.41
27:D1:67:ILE:O	27:D1:68:PRO:O	2.39	0.41
28:D2:22:GLU:HA	28:D2:25:VAL:HG12	2.01	0.41
29:D3:19:GLN:C	29:D3:21:ALA:H	2.24	0.41
31:D5:30:LEU:HD23	31:D5:41:PRO:HA	2.02	0.41
35:DA:811:U:O3'	35:DA:1251:C:H4'	2.21	0.41
35:DA:1287:A:C2'	35:DA:1288:U:H5'	2.50	0.41
35:DA:1417:C:H2'	35:DA:1418:G:C5'	2.51	0.41
35:DA:143(A):C:C2'	35:DA:143(A):C:O2	2.67	0.41
35:DA:1613:G:N1	35:DA:1619:G:C6	2.88	0.41
35:DA:1775:U:C2'	35:DA:1776:G:O5'	2.69	0.41
35:DA:1952:A:C6	35:DA:1953:A:C6	3.09	0.41
35:DA:2199:A:H3'	35:DA:2200:C:H6	1.84	0.41
35:DA:198:C:H5'	35:DA:2244:U:OP1	2.20	0.41
35:DA:2261:C:O4'	35:DA:2388:A:H1'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2574:G:C4	35:DA:2575:C:C5	3.08	0.41
35:DA:2646:C:H2'	35:DA:2647:U:C6	2.55	0.41
35:DA:2818:G:O2'	35:DA:2819:G:H5'	2.20	0.41
35:DA:304:G:N1	35:DA:314:A:C6	2.88	0.41
35:DA:447:A:C4	35:DA:473:G:C8	3.09	0.41
35:DA:702:G:N3	35:DA:731:C:N3	2.68	0.41
35:DA:803:U:C2'	35:DA:804:A:H5'	2.50	0.41
35:DA:808:G:H2'	35:DA:809:G:C8	2.55	0.41
35:DA:836:G:C5	35:DA:837:C:N4	2.89	0.41
35:DA:920:G:C4	35:DA:921:G:C8	3.08	0.41
35:DA:942:G:H1'	35:DA:1189:A:C2	2.55	0.41
36:DB:16:G:O2'	36:DB:17:C:H5'	2.20	0.41
36:DB:49:C:O2'	36:DB:50:G:H5'	2.20	0.41
38:DD:200:ASP:O	38:DD:203:ASN:HB2	2.19	0.41
38:DD:43:ARG:HB3	38:DD:54:ARG:CB	2.50	0.41
39:DE:137:HIS:HB3	39:DE:138:PRO:CD	2.37	0.41
40:DF:90:PHE:O	40:DF:91:GLY:C	2.58	0.41
41:DG:151:ALA:HB3	41:DG:153:ARG:HH12	1.86	0.41
42:DH:121:ILE:HG22	42:DH:133:VAL:CG1	2.50	0.41
42:DH:41:MET:N	42:DH:41:MET:SD	2.93	0.41
43:DI:49:ALA:O	43:DI:53:ALA:HB3	2.21	0.41
44:DN:63:THR:HG23	44:DN:64:GLY:N	2.35	0.41
45:DO:22:ILE:HA	45:DO:22:ILE:HD13	1.80	0.41
45:DO:76:ALA:HB3	50:DT:75:ILE:CG1	2.50	0.41
46:DP:107:LYS:C	46:DP:109:GLY:N	2.73	0.41
46:DP:57:THR:HB	46:DP:59:LEU:HB3	2.02	0.41
46:DP:57:THR:O	46:DP:61:ARG:NH2	2.54	0.41
47:DQ:68:ILE:HG22	47:DQ:101:ARG:HE	1.84	0.41
47:DQ:79:LEU:O	47:DQ:80:GLU:OE2	2.39	0.41
50:DT:100:TYR:C	50:DT:102:ILE:H	2.24	0.41
50:DT:45:PHE:CE2	50:DT:74:ARG:HB2	2.56	0.41
54:DX:38:GLU:O	54:DX:41:ASN:OD1	2.38	0.41
54:DX:60:ARG:HE	54:DX:74:PRO:CG	2.33	0.41
55:DY:37:VAL:HG11	55:DY:72:VAL:HG21	2.01	0.41
1:AA:105:G:C5	1:AA:106:C:C4	3.09	0.41
1:AA:1089:G:H1	1:AA:1096:C:H42	1.69	0.41
1:AA:1242:C:O5'	21:AU:10:ARG:NH1	2.49	0.41
1:AA:1315:U:O2'	1:AA:1360:A:H1'	2.21	0.41
1:AA:1490:C:O2'	1:AA:1491:G:C5'	2.63	0.41
2:AB:220:ASP:OD1	2:AB:220:ASP:N	2.53	0.41
3:AC:127:ARG:NH1	3:AC:127:ARG:CG	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:154:SER:O	3:AC:165:THR:HA	2.21	0.41
3:AC:22:TRP:CZ3	3:AC:32:LEU:HB3	2.56	0.41
5:AE:56:GLN:NE2	5:AE:60:TYR:HB2	2.35	0.41
6:AF:30:LEU:HD23	6:AF:75:LEU:CD2	2.48	0.41
7:AG:36:LYS:O	7:AG:39:ALA:HB3	2.20	0.41
7:AG:62:PHE:O	7:AG:62:PHE:CG	2.73	0.41
8:AH:60:ARG:NH1	8:AH:60:ARG:HG3	2.35	0.41
9:AI:118:LYS:NZ	9:AI:118:LYS:CB	2.81	0.41
9:AI:116:LYS:O	9:AI:119:ALA:N	2.54	0.41
11:AK:33:THR:CA	11:AK:40:ILE:HG12	2.51	0.41
1:AA:562:C:H1'	12:AL:15:ARG:HD2	2.03	0.41
12:AL:82:VAL:HB	12:AL:106:ASP:OD1	2.21	0.41
15:AO:52:SER:O	15:AO:55:GLY:N	2.53	0.41
16:AP:22:THR:HG22	16:AP:32:TYR:CA	2.48	0.41
16:AP:40:ASP:OD2	16:AP:40:ASP:C	2.59	0.41
18:AR:21:LYS:O	18:AR:25:THR:OG1	2.38	0.41
18:AR:33:ASP:OD2	18:AR:34:TYR:N	2.53	0.41
18:AR:40:LEU:O	18:AR:41:LYS:C	2.59	0.41
23:AW:40:C:H2'	23:AW:41:C:C5	2.56	0.41
1:AA:1498:U:C4	24:AX:17:U:H4'	2.56	0.41
25:AY:154:THR:O	25:AY:158:GLU:N	2.46	0.41
25:AY:29:ARG:NE	25:AY:32:ARG:HH22	2.15	0.41
25:AY:35:PRO:HG2	25:AY:36:ALA:H	1.86	0.41
28:B2:29:LYS:O	28:B2:32:LEU:HG	2.20	0.41
32:B6:35:GLU:O	32:B6:35:GLU:HG3	2.21	0.41
35:BA:83:G:N2	35:BA:103:A:OP2	2.53	0.41
35:BA:1174:A:H5''	35:BA:1175:U:C5'	2.50	0.41
35:BA:117:G:H8	35:BA:117:G:O5'	2.04	0.41
35:BA:1261:C:C2'	35:BA:1262:A:O5'	2.69	0.41
35:BA:1374:G:C6	35:BA:1375:C:C4	3.09	0.41
35:BA:1428:C:C5	35:BA:1569:A:H5''	2.55	0.41
35:BA:1691:C:O2'	35:BA:1692:U:H5'	2.20	0.41
35:BA:1805:U:H1'	38:BD:50:THR:O	2.20	0.41
35:BA:1806:C:C5	35:BA:1807:G:N7	2.89	0.41
35:BA:1812:A:O2'	35:BA:1813:G:H5'	2.21	0.41
35:BA:185:U:O2'	35:BA:186:G:H5'	2.21	0.41
35:BA:213:A:C2'	35:BA:214:G:H5'	2.50	0.41
35:BA:2241:A:N1	35:BA:2242:G:C6	2.88	0.41
35:BA:2473:U:O2	35:BA:2473:U:H2'	2.20	0.41
35:BA:2522:U:C2'	35:BA:2523:G:H5''	2.50	0.41
35:BA:2711:A:C8	35:BA:2714:G:O4'	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1661:G:H5'	35:BA:2712:U:O4	2.20	0.41
35:BA:271(S):G:H2'	35:BA:271(T):C:O4'	2.21	0.41
35:BA:2751:G:N3	35:BA:2751:G:H2'	2.36	0.41
35:BA:2783:G:N2	35:BA:2784:C:C2	2.89	0.41
35:BA:593:G:H2'	35:BA:594:U:C6	2.55	0.41
35:BA:596:G:C2	35:BA:662:G:C2	3.08	0.41
35:BA:669:G:O2'	35:BA:670:A:OP1	2.29	0.41
35:BA:696:G:H2'	35:BA:697:C:H5'	2.02	0.41
35:BA:790:C:O2'	35:BA:791:C:C5'	2.69	0.41
35:BA:856:C:H5''	35:BA:856:C:H6	1.84	0.41
35:BA:995:C:C2	51:BU:57:PHE:HE2	2.38	0.41
35:BA:998:C:OP2	51:BU:93:LYS:HE2	2.21	0.41
38:BD:61:LEU:HD12	38:BD:62:TYR:N	2.35	0.41
38:BD:62:TYR:OH	38:BD:64:ILE:HD12	2.20	0.41
39:BE:30:PRO:CD	39:BE:180:ASN:HD21	2.23	0.41
39:BE:44:TYR:CE2	39:BE:46:ALA:HB2	2.56	0.41
39:BE:52:LEU:N	39:BE:74:PRO:CB	2.83	0.41
40:BF:110:LEU:HD23	40:BF:110:LEU:C	2.40	0.41
40:BF:139:PHE:O	40:BF:140:LEU:C	2.59	0.41
40:BF:20:LEU:HB2	40:BF:24:LEU:HD21	2.02	0.41
40:BF:47:GLY:HA3	40:BF:95:ARG:O	2.20	0.41
41:BG:115:ARG:HH22	41:BG:136:ARG:CG	2.33	0.41
41:BG:124:SER:HB2	41:BG:131:TYR:CE1	2.56	0.41
42:BH:96:ALA:CB	42:BH:105:LEU:HB3	2.48	0.41
43:BI:122:GLU:O	43:BI:126:TYR:OH	2.39	0.41
43:BI:81:VAL:O	43:BI:83:ALA:N	2.53	0.41
44:BN:2:LYS:HZ1	52:BV:12:TYR:HB3	1.85	0.41
44:BN:55:VAL:CG1	44:BN:56:ASN:N	2.84	0.41
46:BP:57:THR:C	46:BP:59:LEU:N	2.74	0.41
47:BQ:29:PHE:N	47:BQ:105:GLU:OE2	2.44	0.41
47:BQ:12:GLN:HE21	47:BQ:73:PRO:HD3	1.85	0.41
49:BS:53:SER:O	49:BS:55:ALA:N	2.54	0.41
50:BT:109:GLU:C	50:BT:112:ARG:HG3	2.40	0.41
50:BT:33:LYS:HB2	50:BT:41:ARG:HB3	2.03	0.41
50:BT:61:PHE:CE1	50:BT:76:PHE:HD1	2.38	0.41
53:BW:59:VAL:C	53:BW:61:ASN:H	2.23	0.41
53:BW:27:LYS:O	53:BW:71:VAL:HG23	2.21	0.41
54:BX:36:LYS:HZ2	54:BX:39:ILE:HA	1.81	0.41
56:BZ:125:LEU:O	56:BZ:165:VAL:HG23	2.20	0.41
1:CA:1002:G:H2'	1:CA:1003:G:C8	2.56	0.41
1:CA:1151:A:C4	1:CA:1152:A:N7	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1232:U:H5'	9:CI:126:SER:OG	2.19	0.41
1:CA:329:A:C2	1:CA:332:G:C5	3.09	0.41
1:CA:620:C:O2'	1:CA:621:A:H5'	2.20	0.41
1:CA:751:U:H2'	1:CA:752:G:O4'	2.20	0.41
1:CA:769:G:C4	1:CA:770:C:C5	3.09	0.41
1:CA:824:C:H2'	1:CA:825:G:H8	1.85	0.41
1:CA:874:G:O2'	1:CA:875:C:H5'	2.20	0.41
1:CA:935:A:H61	1:CA:1380:U:H3	1.69	0.41
2:CB:187:LEU:HD22	2:CB:188:ALA:O	2.21	0.41
3:CC:134:ILE:CG2	3:CC:151:VAL:HB	2.51	0.41
3:CC:6:HIS:HA	3:CC:7:PRO:HD2	1.91	0.41
4:CD:21:LEU:HD22	4:CD:115:ARG:CG	2.51	0.41
4:CD:129:ASN:HB2	4:CD:131:ARG:HH22	1.85	0.41
4:CD:160:GLN:O	4:CD:163:GLU:HB3	2.21	0.41
5:CE:128:PRO:O	5:CE:131:ILE:HG12	2.21	0.41
5:CE:78:HIS:HB2	5:CE:79:GLU:H	1.69	0.41
6:CF:100:ASN:HA	6:CF:100:ASN:HD22	1.54	0.41
8:CH:54:ASP:C	8:CH:56:LYS:H	2.23	0.41
10:CJ:64:GLU:O	10:CJ:64:GLU:HG3	2.21	0.41
12:CL:44:THR:HA	12:CL:45:PRO:HD3	1.85	0.41
13:CM:19:LEU:O	13:CM:22:ILE:HB	2.20	0.41
16:CP:53:VAL:O	16:CP:54:GLU:C	2.58	0.41
18:CR:40:LEU:O	18:CR:41:LYS:C	2.58	0.41
25:CY:23:HIS:O	25:CY:24:ASN:C	2.58	0.41
25:CY:45:TYR:CD2	25:CY:78:ALA:HB2	2.55	0.41
28:D2:60:LEU:CG	28:D2:61:LEU:H	2.10	0.41
33:D7:24:THR:O	33:D7:28:ARG:HG3	2.20	0.41
35:DA:1450:G:H2'	35:DA:1450(A):C:C6	2.55	0.41
35:DA:1528(A):A:H3'	35:DA:1529:G:C5'	2.45	0.41
35:DA:1555:G:H2'	35:DA:1556:C:H6	1.82	0.41
35:DA:1906:G:C8	35:DA:1929:G:C4	3.09	0.41
35:DA:1986:A:H2'	35:DA:1987:G:H5''	2.02	0.41
35:DA:2053:G:N2	35:DA:2054:A:C4	2.89	0.41
35:DA:2197:U:H3	35:DA:2225:A:H62	1.69	0.41
35:DA:2253:G:C5	35:DA:2254:C:C4	3.09	0.41
35:DA:2298:A:C2	35:DA:2321:G:C4	3.09	0.41
35:DA:271(V):G:H2'	35:DA:271(W):G:C8	2.56	0.41
35:DA:2785:C:H6	35:DA:2785:C:O5'	2.04	0.41
35:DA:2811:G:N2	35:DA:2891:G:H1'	2.36	0.41
35:DA:385:C:C2'	35:DA:386:G:OP2	2.68	0.41
35:DA:485:C:C2	35:DA:496:G:N2	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:832:G:OP1	46:DP:40:SER:HB3	2.21	0.41
35:DA:8:A:OP1	44:DN:51:PHE:HE2	2.04	0.41
35:DA:976:C:H42	35:DA:988:A:H2	1.68	0.41
35:DA:998:C:OP2	51:DU:93:LYS:CE	2.68	0.41
36:DB:54:G:O2'	36:DB:55:U:H5'	2.21	0.41
38:DD:138:VAL:HG21	38:DD:166:GLN:O	2.21	0.41
38:DD:80:ALA:O	38:DD:81:ALA:HB2	2.20	0.41
39:DE:13:ARG:HA	39:DE:21:VAL:O	2.21	0.41
39:DE:52:LEU:O	39:DE:53:PRO:O	2.38	0.41
39:DE:61:ARG:CG	39:DE:62:PRO:N	2.84	0.41
40:DF:25:PRO:HB3	40:DF:119:ARG:CG	2.51	0.41
40:DF:53:THR:CG2	40:DF:56:GLU:OE1	2.69	0.41
41:DG:116:ASP:CG	41:DG:117:PHE:N	2.74	0.41
42:DH:111:HIS:HA	42:DH:112:PRO:HD2	1.95	0.41
42:DH:122:THR:CB	42:DH:134:SER:HB2	2.46	0.41
44:DN:61:ARG:HG3	44:DN:61:ARG:NH1	2.35	0.41
46:DP:30:THR:O	46:DP:33:ARG:N	2.49	0.41
46:DP:70:GLN:CG	46:DP:71:VAL:N	2.83	0.41
47:DQ:54:MET:CG	47:DQ:64:ILE:HD13	2.50	0.41
47:DQ:68:ILE:HG12	47:DQ:68:ILE:O	2.21	0.41
47:DQ:89:ASN:O	47:DQ:91:GLU:N	2.53	0.41
48:DR:26:LYS:HG2	48:DR:70:LEU:HD22	2.02	0.41
48:DR:28:LEU:HD12	48:DR:29:LEU:CD1	2.51	0.41
48:DR:67:LEU:HA	48:DR:67:LEU:HD13	1.86	0.41
48:DR:56:LYS:HD2	48:DR:88:ARG:HA	2.02	0.41
49:DS:74:ALA:O	49:DS:78:LEU:HG	2.21	0.41
36:DB:47:C:O2'	49:DS:93:LYS:HG3	2.21	0.41
50:DT:72:VAL:CG1	50:DT:73:GLU:N	2.83	0.41
51:DU:36:ARG:O	51:DU:37:GLU:C	2.57	0.41
53:DW:64:MET:HG2	53:DW:109:GLU:OE2	2.20	0.41
55:DY:28:LYS:O	55:DY:29:GLU:C	2.59	0.41
56:DZ:134:PRO:HG3	56:DZ:161:VAL:HG21	2.02	0.41
56:DZ:43:GLU:O	56:DZ:47:VAL:HG23	2.20	0.41
56:DZ:76:LEU:CD2	56:DZ:76:LEU:H	2.32	0.41
1:AA:1421:G:C6	1:AA:1422:G:C5	3.09	0.41
1:AA:1501:C:N4	1:AA:1504:G:C2	2.88	0.41
1:AA:258:G:H2'	1:AA:259:G:H8	1.86	0.41
1:AA:27:G:H2'	1:AA:28:G:O4'	2.20	0.41
1:AA:779:C:C2'	1:AA:780:A:H5'	2.50	0.41
3:AC:119:ARG:HH21	3:AC:140:ARG:NE	2.19	0.41
4:AD:121:VAL:CA	4:AD:126:ILE:HD12	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:148:VAL:HG23	4:AD:181:MET:O	2.20	0.41
4:AD:150:GLU:HA	4:AD:153:ARG:HD3	2.03	0.41
4:AD:66:ARG:O	4:AD:69:GLY:N	2.54	0.41
4:AD:97:LEU:O	4:AD:97:LEU:HD22	2.20	0.41
5:AE:102:ALA:H	5:AE:107:ARG:NH2	2.15	0.41
7:AG:42:ILE:HA	7:AG:45:ASP:CB	2.49	0.41
7:AG:67:GLU:CD	7:AG:67:GLU:H	2.24	0.41
8:AH:127:LEU:HD22	8:AH:127:LEU:HA	1.85	0.41
8:AH:26:VAL:CG2	8:AH:32:LYS:NZ	2.76	0.41
9:AI:102:LEU:C	9:AI:102:LEU:HD23	2.41	0.41
1:AA:1179:A:O2'	9:AI:103:THR:HG23	2.20	0.41
9:AI:5:TYR:HE2	9:AI:16:ARG:HB3	1.86	0.41
9:AI:28:VAL:CG1	9:AI:29:ASN:H	2.32	0.41
1:AA:1060:C:H4'	10:AJ:52:GLY:CA	2.51	0.41
11:AK:108:ILE:HD12	11:AK:108:ILE:N	2.35	0.41
11:AK:22:HIS:HB3	11:AK:29:ILE:CG1	2.36	0.41
12:AL:47:LYS:HG2	12:AL:48:PRO:HD3	2.01	0.41
1:AA:995:C:H1'	14:AN:4:LYS:HE3	2.02	0.41
20:AT:12:ALA:C	20:AT:14:LYS:N	2.74	0.41
20:AT:14:LYS:CA	20:AT:17:ARG:HH21	2.31	0.41
23:AW:47:G:H2'	23:AW:48:U:H5'	2.03	0.41
27:B1:71:TYR:HA	27:B1:74:VAL:HG21	2.00	0.41
34:B8:58:ILE:HA	34:B8:61:LEU:HD21	2.03	0.41
35:BA:1013:C:O2'	35:BA:1014:U:H5'	2.20	0.41
35:BA:84:A:N6	35:BA:102:G:O2'	2.54	0.41
35:BA:108:U:C2	35:BA:109:G:N7	2.89	0.41
35:BA:115:C:H2'	35:BA:116:C:H6	1.86	0.41
35:BA:1181:C:H2'	35:BA:1182:A:H8	1.86	0.41
35:BA:1389:G:H2'	35:BA:1390:U:C6	2.56	0.41
35:BA:1474:C:H3'	35:BA:1475:G:H8	1.84	0.41
35:BA:149:A:H2'	35:BA:150:C:H6	1.82	0.41
35:BA:1844:C:H5''	38:BD:258:LYS:HG3	2.02	0.41
35:BA:1992:G:HO2'	35:BA:1993:U:P	2.44	0.41
35:BA:2061:G:O4'	35:BA:2503:A:C5	2.74	0.41
35:BA:2186:G:H3'	35:BA:2187:G:H5''	2.02	0.41
35:BA:2206:G:N3	35:BA:2207:G:H5'	2.34	0.41
35:BA:2234:G:H2'	35:BA:2235:G:H8	1.86	0.41
35:BA:2288:A:H2	35:BA:2325:G:C8	2.39	0.41
35:BA:2472:G:H2'	35:BA:2529:G:N2	2.35	0.41
35:BA:2543:G:O2'	35:BA:2544:G:H5'	2.21	0.41
35:BA:2562:U:O2'	35:BA:2563:U:H5'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:2602:A:H4'	35:BA:2603:G:O5'	2.21	0.41
35:BA:2673:G:H2'	35:BA:2674:G:C8	2.50	0.41
35:BA:2757:A:N1	42:BH:67:LEU:HD13	2.36	0.41
35:BA:814:C:H3'	57:BA:3399:MG:MG	1.45	0.41
35:BA:359:A:H2'	35:BA:360:G:O4'	2.21	0.41
35:BA:372:G:O2'	35:BA:373:U:OP2	2.38	0.41
35:BA:511:U:O4	35:BA:512:G:C2	2.73	0.41
35:BA:589:C:O3'	40:BF:95:ARG:NH1	2.53	0.41
38:BD:178:PRO:C	38:BD:180:GLY:H	2.24	0.41
38:BD:25:THR:HB	38:BD:82:ILE:N	2.24	0.41
39:BE:101:ARG:HH11	39:BE:169:ASN:HD22	1.68	0.41
39:BE:137:HIS:HB3	39:BE:138:PRO:CD	2.41	0.41
39:BE:13:ARG:HA	39:BE:21:VAL:O	2.19	0.41
39:BE:67:PHE:HA	39:BE:67:PHE:HD2	1.76	0.41
40:BF:148:LEU:HD13	40:BF:154:VAL:HG21	2.02	0.41
40:BF:24:LEU:HD22	40:BF:24:LEU:N	2.36	0.41
41:BG:17:PRO:HA	41:BG:20:ILE:CG1	2.46	0.41
41:BG:34:LEU:HB3	41:BG:99:MET:SD	2.60	0.41
43:BI:10:GLU:C	43:BI:12:LEU:H	2.24	0.41
35:BA:6:A:C2'	44:BN:130:HIS:HB2	2.50	0.41
44:BN:67:LEU:HB3	44:BN:88:GLU:CG	2.51	0.41
45:BO:61:VAL:O	45:BO:61:VAL:HG13	2.21	0.41
46:BP:131:SER:CB	46:BP:134:ALA:HB3	2.46	0.41
46:BP:138:LEU:HD22	46:BP:142:GLY:HA3	2.02	0.41
35:BA:662:G:P	46:BP:18:ARG:HD2	2.61	0.41
35:BA:245:G:H5'	46:BP:70:GLN:H	1.86	0.41
48:BR:18:LEU:HD13	48:BR:19:ALA:CA	2.50	0.41
50:BT:92:GLY:O	50:BT:94:ALA:N	2.54	0.41
51:BU:106:PHE:O	51:BU:107:ALA:C	2.58	0.41
51:BU:31:SER:HB3	51:BU:34:LYS:CG	2.50	0.41
51:BU:92:ARG:CZ	52:BV:11:GLN:H	2.33	0.41
52:BV:88:ARG:CG	52:BV:88:ARG:HH11	2.33	0.41
52:BV:73:SER:OG	52:BV:89:GLN:O	2.30	0.41
53:BW:1:MET:O	53:BW:2:GLU:HB3	2.20	0.41
54:BX:60:ARG:HG2	54:BX:73:ARG:N	2.36	0.41
55:BY:87:LYS:HG2	55:BY:88:LYS:H	1.86	0.41
55:BY:90:LEU:N	55:BY:90:LEU:HD23	2.36	0.41
1:CA:1184:G:C2	1:CA:1185:G:C5	3.08	0.41
1:CA:1342:C:O2'	1:CA:1343:G:H5'	2.21	0.41
1:CA:286:G:H2'	1:CA:287:U:H6	1.85	0.41
1:CA:657:G:C2	1:CA:658:G:C8	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:792:A:N3	1:CA:794:A:C5	2.89	0.41
1:CA:857:C:O2	1:CA:858:G:H1'	2.20	0.41
2:CB:100:GLY:CA	2:CB:103:THR:HB	2.50	0.41
2:CB:77:ALA:HA	2:CB:208:ILE:HG12	2.02	0.41
2:CB:212:GLN:O	2:CB:213:LEU:C	2.58	0.41
1:CA:1107:C:OP1	3:CC:174:PRO:HD3	2.20	0.41
3:CC:150:LYS:CB	3:CC:201:TYR:HB2	2.38	0.41
4:CD:18:LYS:HE3	4:CD:31:CYS:CB	2.51	0.41
4:CD:33:MET:HE2	4:CD:33:MET:HA	2.03	0.41
6:CF:16:GLN:O	6:CF:17:SER:C	2.58	0.41
5:CE:78:HIS:CD2	8:CH:104:ARG:NE	2.89	0.41
8:CH:95:VAL:HA	8:CH:99:GLU:OE1	2.21	0.41
9:CI:47:LEU:CB	9:CI:50:LEU:HD12	2.48	0.41
9:CI:63:ILE:CD1	9:CI:63:ILE:N	2.80	0.41
10:CJ:6:ILE:CD1	10:CJ:72:VAL:HB	2.51	0.41
12:CL:19:ARG:HD2	12:CL:19:ARG:H	1.86	0.41
15:CO:29:VAL:O	15:CO:30:ALA:C	2.58	0.41
16:CP:4:ILE:HA	16:CP:21:VAL:HA	2.02	0.41
1:CA:375:U:O3'	16:CP:6:LEU:HD22	2.21	0.41
17:CQ:92:ARG:O	17:CQ:93:GLN:C	2.59	0.41
20:CT:40:ALA:HB2	20:CT:55:ILE:CG2	2.50	0.41
23:CW:34:U:O2	23:CW:37:U:C5	2.74	0.41
25:CY:109:GLU:C	25:CY:112:LYS:HB3	2.40	0.41
25:CY:32:ARG:HB3	25:CY:103:ILE:CD1	2.48	0.41
26:D0:1:MET:O	26:D0:2:ALA:CB	2.68	0.41
26:D0:32:ARG:HB2	26:D0:35:ASN:ND2	2.36	0.41
26:D0:32:ARG:HB3	26:D0:33:ALA:H	1.70	0.41
31:D5:17:ASP:HA	31:D5:20:ARG:CG	2.51	0.41
33:D7:25:PRO:HA	33:D7:28:ARG:CZ	2.50	0.41
35:DA:139(A):G:H5'	35:DA:140:G:OP2	2.20	0.41
35:DA:1449:A:C6	35:DA:1450:G:H1'	2.55	0.41
35:DA:1477:A:C2	35:DA:1515:G:C2	3.09	0.41
35:DA:1271:G:N2	35:DA:1617:C:O4'	2.54	0.41
35:DA:1765:C:O5'	35:DA:1765:C:H6	2.04	0.41
35:DA:1767:C:O2'	35:DA:1768:U:H5'	2.21	0.41
35:DA:1861:G:H2'	35:DA:1862:G:C8	2.55	0.41
35:DA:2016:U:O5'	35:DA:2016:U:H6	2.03	0.41
35:DA:2027:G:C5	35:DA:2028:U:C5	3.09	0.41
35:DA:2068:U:O4	35:DA:2430:A:C2	2.74	0.41
35:DA:2127:G:H1'	35:DA:2128:C:C4'	2.47	0.41
35:DA:2219:G:C6	35:DA:2220:G:N7	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:2346:A:H5'	35:DA:2383:G:O4'	2.21	0.41
35:DA:2251:G:C2	35:DA:2450:A:H1'	2.56	0.41
35:DA:2657:A:C5	35:DA:2658:C:C5	3.09	0.41
35:DA:2721:A:H2'	35:DA:2722:G:O4'	2.21	0.41
35:DA:2751:G:H2'	35:DA:2751:G:N3	2.36	0.41
35:DA:27:G:H1'	35:DA:513:A:N6	2.35	0.41
35:DA:2805:G:C2	35:DA:2807:G:C5	3.09	0.41
35:DA:610:G:C6	35:DA:611:C:N4	2.87	0.41
35:DA:645:C:O2	35:DA:645:C:H2'	2.21	0.41
35:DA:677:A:N1	35:DA:802:A:C6	2.89	0.41
35:DA:695:G:C2	35:DA:696:G:C8	3.09	0.41
35:DA:761:A:O5'	35:DA:761:A:H8	2.04	0.41
35:DA:900:A:N7	35:DA:901:A:N7	2.68	0.41
35:DA:980:A:C6	35:DA:981:A:C2	3.09	0.41
38:DD:142:VAL:HG23	38:DD:192:THR:O	2.20	0.41
38:DD:142:VAL:CG2	38:DD:143:HIS:N	2.84	0.41
38:DD:35:LYS:HG2	38:DD:64:ILE:CG2	2.51	0.41
35:DA:2724:C:OP1	39:DE:118:LYS:HE3	2.20	0.41
39:DE:120:TRP:C	39:DE:122:PHE:N	2.72	0.41
35:DA:2572:A:C6	39:DE:144:ARG:NH2	2.89	0.41
40:DF:170:LEU:HD23	40:DF:173:VAL:HG21	2.01	0.41
40:DF:26:ALA:HB1	40:DF:27:GLU:OE1	2.21	0.41
40:DF:7:TYR:CD1	40:DF:196:LEU:HD11	2.56	0.41
40:DF:81:PRO:HG2	40:DF:82:ILE:HG23	2.03	0.41
41:DG:108:ASN:O	41:DG:112:PRO:HD2	2.20	0.41
42:DH:12:PRO:O	42:DH:13:LYS:CB	2.68	0.41
43:DI:86:THR:O	43:DI:122:GLU:OE2	2.39	0.41
43:DI:132:PRO:O	43:DI:133:HIS:C	2.59	0.41
43:DI:130:TYR:O	43:DI:135:GLU:HG2	2.21	0.41
43:DI:44:LEU:C	43:DI:44:LEU:HD23	2.41	0.41
44:DN:18:ALA:C	44:DN:20:GLY:N	2.74	0.41
45:DO:47:ILE:HG23	45:DO:48:PRO:N	2.36	0.41
45:DO:88:ASN:O	45:DO:89:ASN:C	2.59	0.41
47:DQ:87:LYS:CG	47:DQ:87:LYS:O	2.55	0.41
31:D5:44:THR:CG2	48:DR:101:ALA:N	2.84	0.41
48:DR:15:SER:O	48:DR:16:HIS:O	2.39	0.41
48:DR:72:ASP:HB3	48:DR:75:LEU:HB2	2.03	0.41
50:DT:10:VAL:N	50:DT:12:SER:OG	2.53	0.41
51:DU:7:GLY:C	51:DU:8:VAL:CG2	2.89	0.41
52:DV:4:ILE:CD1	52:DV:40:LEU:HD11	2.49	0.41
52:DV:63:GLY:O	52:DV:64:HIS:CB	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:DX:81:VAL:C	54:DX:82:GLN:O	2.58	0.41
54:DX:89:ILE:HA	54:DX:92:LEU:HB2	2.01	0.41
56:DZ:7:ALA:HB1	56:DZ:37:VAL:HG21	2.02	0.41
56:DZ:72:ARG:CG	56:DZ:89:PHE:HB2	2.43	0.41
1:AA:113:G:H2'	1:AA:114:U:C6	2.55	0.41
1:AA:1226:C:H2'	13:AM:103:THR:CB	2.51	0.41
1:AA:1421:G:O2'	1:AA:1422:G:H5'	2.21	0.41
1:AA:1442:G:C4	1:AA:1442(B):A:N1	2.89	0.41
1:AA:66:G:O4'	1:AA:173:U:C4	2.74	0.41
1:AA:18:C:H2'	1:AA:19:C:H6	1.86	0.41
1:AA:253:U:O2'	1:AA:254:G:H5'	2.20	0.41
1:AA:451:A:C1'	1:AA:452:A:N7	2.84	0.41
1:AA:52:G:C6	1:AA:360:A:C2	3.09	0.41
1:AA:602:A:O2'	1:AA:603:U:H5'	2.20	0.41
1:AA:656:C:O2'	1:AA:657:G:H5'	2.20	0.41
1:AA:769:G:H2'	1:AA:770:C:C6	2.52	0.41
2:AB:115:LEU:CB	2:AB:145:LEU:HD11	2.50	0.41
2:AB:69:LEU:HB3	2:AB:162:ILE:HB	2.03	0.41
2:AB:70:PHE:O	2:AB:92:TYR:CB	2.66	0.41
2:AB:83:MET:HB2	2:AB:84:GLU:OE2	2.21	0.41
3:AC:95:THR:HB	3:AC:99:VAL:CG1	2.51	0.41
4:AD:158:ILE:HG22	4:AD:159:ARG:N	2.35	0.41
5:AE:69:VAL:HG11	5:AE:113:ALA:HB1	2.03	0.41
7:AG:64:GLN:HG2	7:AG:128:ALA:HB1	2.03	0.41
7:AG:23:VAL:HG12	7:AG:27:ILE:CD1	2.51	0.41
8:AH:3:THR:O	8:AH:5:PRO:HD3	2.21	0.41
8:AH:95:VAL:HA	8:AH:99:GLU:OE1	2.21	0.41
8:AH:9:MET:C	8:AH:11:THR:N	2.73	0.41
9:AI:28:VAL:HG22	9:AI:63:ILE:C	2.41	0.41
10:AJ:45:ARG:NH1	14:AN:36:PHE:CE2	2.89	0.41
13:AM:19:LEU:HA	13:AM:22:ILE:CD1	2.38	0.41
13:AM:48:LEU:HD21	13:AM:53:VAL:CG2	2.51	0.41
16:AP:65:GLN:HA	16:AP:66:PRO:HD2	1.93	0.41
17:AQ:45:HIS:HE2	17:AQ:47:PRO:HB3	1.84	0.41
20:AT:26:ASN:HD22	20:AT:27:LYS:N	2.19	0.41
25:AY:164:ILE:CG2	25:AY:165:THR:N	2.80	0.41
25:AY:48:ALA:O	25:AY:50:VAL:N	2.54	0.41
25:AY:61:PRO:HG3	25:AY:67:VAL:CG1	2.50	0.41
27:B1:13:ILE:HD13	27:B1:66:HIS:CE1	2.56	0.41
27:B1:75:GLU:O	27:B1:76:ARG:CZ	2.69	0.41
34:B8:2:PRO:HA	35:BA:591:C:O2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:B8:37:SER:HB2	34:B8:38:GLY:H	1.67	0.41
35:BA:1109:C:N4	35:BA:1110:G:C2	2.89	0.41
35:BA:1208:C:C4	35:BA:1209:G:N7	2.89	0.41
35:BA:1272:A:C2	35:BA:1618:A:C5	3.09	0.41
35:BA:1353:A:H4'	38:BD:38:LYS:HZ3	1.86	0.41
35:BA:769:G:H5'	35:BA:1379:A:H61	1.86	0.41
35:BA:1441:G:H4'	35:BA:1627:G:O3'	2.21	0.41
35:BA:1692:U:H2'	35:BA:1694:C:C5	2.55	0.41
35:BA:1757:U:H2'	35:BA:1758:G:OP1	2.21	0.41
35:BA:1785:A:N7	35:BA:1787:A:C5	2.89	0.41
35:BA:1925:C:H2'	35:BA:1926:U:C5'	2.50	0.41
35:BA:2036:C:C6	35:BA:2036:C:C4'	3.04	0.41
35:BA:2253:G:C5	35:BA:2254:C:C4	3.08	0.41
26:B0:20:ARG:HB3	35:BA:2270:G:O3'	2.21	0.41
35:BA:2261:C:C2	35:BA:2280:G:C2	3.08	0.41
35:BA:2517:C:C6	35:BA:2542:A:C2	3.09	0.41
35:BA:271(W):G:C5'	35:BA:271(X):G:OP2	2.69	0.41
35:BA:26:G:C6	35:BA:27:G:C2	3.09	0.41
35:BA:348:G:H2'	35:BA:349:G:C8	2.56	0.41
35:BA:942:G:H2'	35:BA:943:U:O4'	2.20	0.41
35:BA:1567:A:N6	38:BD:21:PHE:CE2	2.89	0.41
38:BD:70:TRP:CZ2	38:BD:150:LYS:HA	2.55	0.41
38:BD:8:PRO:HB3	38:BD:14:ARG:HA	2.03	0.41
42:BH:94:TYR:CE1	42:BH:160:LYS:HD2	2.55	0.41
42:BH:41:MET:HE3	42:BH:54:ARG:HA	2.03	0.41
42:BH:35:VAL:HG21	42:BH:72:ILE:HD13	2.02	0.41
42:BH:83:TYR:CD1	42:BH:84:SER:N	2.89	0.41
44:BN:110:GLY:HA2	44:BN:114:ARG:HH21	1.81	0.41
45:BO:13:ASN:ND2	45:BO:96:THR:OG1	2.53	0.41
48:BR:78:LYS:O	48:BR:82:GLU:HB3	2.20	0.41
49:BS:34:HIS:NE2	49:BS:54:LEU:HB2	2.36	0.41
53:BW:44:ALA:O	53:BW:45:TYR:C	2.58	0.41
55:BY:87:LYS:C	55:BY:89:PHE:H	2.23	0.41
55:BY:87:LYS:HG2	55:BY:88:LYS:N	2.35	0.41
56:BZ:91:LEU:HB3	56:BZ:130:PRO:CG	2.51	0.41
1:CA:1019:C:O2'	1:CA:1020:U:H5'	2.21	0.41
1:CA:1226:C:OP1	19:CS:81:ARG:CZ	2.69	0.41
1:CA:1249:C:H6	1:CA:1249:C:H5'	1.85	0.41
1:CA:1517:G:O6	1:CA:1518:A:N1	2.53	0.41
1:CA:1529:G:OP2	1:CA:1529:G:H3'	2.21	0.41
1:CA:37:U:O2'	1:CA:38:G:H5'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:80:G:H3'	1:CA:81:U:H5'	2.03	0.41
1:CA:810:C:H2'	1:CA:811:C:O4'	2.21	0.41
1:CA:885:G:H1	1:CA:912:C:N4	2.18	0.41
1:CA:919:A:C2'	1:CA:920:U:O5'	2.69	0.41
1:CA:947:G:H2'	1:CA:948:C:C6	2.56	0.41
1:CA:974:A:H8	1:CA:974:A:OP1	2.04	0.41
2:CB:185:ILE:HG22	2:CB:199:TYR:CB	2.45	0.41
2:CB:43:ASP:OD2	2:CB:46:LYS:N	2.46	0.41
3:CC:128:PHE:O	3:CC:130:VAL:N	2.54	0.41
3:CC:89:GLU:O	3:CC:93:LYS:N	2.49	0.41
4:CD:53:ASP:HB3	4:CD:57:ARG:NH2	2.36	0.41
5:CE:11:ILE:HD12	5:CE:31:LEU:CD2	2.48	0.41
1:CA:18:C:C5'	5:CE:127:ASN:ND2	2.84	0.41
6:CF:18:GLN:CA	6:CF:21:LEU:HD23	2.46	0.41
8:CH:120:THR:C	8:CH:122:ARG:N	2.74	0.41
8:CH:20:TYR:HA	8:CH:65:TYR:CZ	2.55	0.41
8:CH:54:ASP:C	8:CH:56:LYS:N	2.74	0.41
9:CI:83:ARG:HA	9:CI:86:VAL:HG12	2.03	0.41
10:CJ:22:LYS:HE3	10:CJ:23:ILE:HG13	2.03	0.41
11:CK:71:LYS:O	11:CK:73:MET:N	2.54	0.41
13:CM:113:PRO:O	13:CM:114:ARG:C	2.58	0.41
17:CQ:45:HIS:HB2	17:CQ:69:LYS:CE	2.48	0.41
18:CR:71:LYS:C	18:CR:74:ARG:HB2	2.41	0.41
19:CS:48:THR:HG22	19:CS:61:TYR:CB	2.50	0.41
19:CS:70:LYS:NZ	19:CS:70:LYS:CB	2.77	0.41
20:CT:14:LYS:CA	20:CT:17:ARG:HH21	2.33	0.41
20:CT:48:LYS:O	20:CT:49:ALA:CB	2.69	0.41
25:CY:104:PRO:HA	25:CY:105:PRO:HD2	1.81	0.41
25:CY:152:ASP:O	25:CY:156:ARG:HG3	2.21	0.41
25:CY:83:ILE:C	25:CY:85:ASP:N	2.72	0.41
26:D0:25:ARG:CB	26:D0:37:LEU:HD23	2.49	0.41
26:D0:40:GLN:HG3	26:D0:57:PHE:HB3	2.03	0.41
27:D1:78:LYS:HE2	35:DA:271(R):G:O4'	2.21	0.41
28:D2:37:PHE:CD2	28:D2:40:SER:HA	2.56	0.41
31:D5:27:PRO:O	31:D5:28:PRO:C	2.59	0.41
31:D5:50:GLY:HA3	31:D5:56:LYS:HB3	2.03	0.41
33:D7:25:PRO:HG2	33:D7:26:GLY:N	2.31	0.41
34:D8:59:LYS:O	34:D8:61:LEU:N	2.50	0.41
35:DA:1488:G:H2'	35:DA:1488:G:N3	2.36	0.41
35:DA:1779:U:P	35:DA:1784:A:H61	2.43	0.41
35:DA:1826:G:H4'	38:DD:242:ARG:CZ	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1663:C:N3	35:DA:1992:G:O6	2.54	0.41
35:DA:2065:C:H1'	35:DA:2449:U:N3	2.34	0.41
35:DA:2073:C:O2'	35:DA:2074:U:H5'	2.20	0.41
35:DA:2171:A:N3	35:DA:2171:A:H2'	2.36	0.41
35:DA:2297:C:H2'	35:DA:2298:A:C5'	2.36	0.41
35:DA:2571:C:O2	35:DA:2571:C:C2'	2.67	0.41
35:DA:2753:A:H2	35:DA:2754:U:N3	2.18	0.41
35:DA:363(F):A:O2'	35:DA:364:C:C5	2.74	0.41
35:DA:481:G:O2'	35:DA:482:A:P	2.79	0.41
35:DA:532:A:C8	35:DA:2021:C:C5	3.09	0.41
35:DA:535:C:C4	35:DA:536:A:N7	2.89	0.41
35:DA:562:U:O2'	35:DA:563:G:C5'	2.69	0.41
35:DA:589:C:O3'	40:DF:95:ARG:NH1	2.54	0.41
35:DA:802:A:C6	35:DA:803:U:C4	3.09	0.41
35:DA:968:G:C2	35:DA:969:U:C2	3.09	0.41
36:DB:11:C:H3'	36:DB:12:C:C5	2.56	0.41
38:DD:166:GLN:CA	38:DD:166:GLN:NE2	2.83	0.41
38:DD:53:PHE:C	38:DD:54:ARG:HG2	2.38	0.41
39:DE:117:MET:HE1	39:DE:124:GLY:HA3	2.02	0.41
39:DE:35:GLN:HE22	39:DE:37:ARG:HH21	1.68	0.41
39:DE:56:PRO:O	39:DE:57:LYS:C	2.59	0.41
40:DF:110:LEU:C	40:DF:110:LEU:HD23	2.40	0.41
40:DF:171:PRO:C	40:DF:173:VAL:H	2.24	0.41
40:DF:82:ILE:H	40:DF:82:ILE:HG23	1.58	0.41
41:DG:121:ASN:OD1	41:DG:123:ASN:OD1	2.38	0.41
41:DG:181:ARG:O	41:DG:182:LYS:OXT	2.39	0.41
41:DG:45:GLU:HA	41:DG:47:LYS:HD3	2.03	0.41
42:DH:58:GLU:O	42:DH:61:HIS:HB2	2.20	0.41
44:DN:87:LEU:O	44:DN:90:MET:HB2	2.21	0.41
44:DN:96:GLU:OE2	44:DN:96:GLU:N	2.51	0.41
45:DO:60:ALA:HA	45:DO:87:ILE:H	1.86	0.41
46:DP:101:VAL:CB	46:DP:107:LYS:HA	2.50	0.41
46:DP:13:ASN:ND2	46:DP:13:ASN:N	2.69	0.41
46:DP:48:PRO:CG	46:DP:49:ARG:H	2.34	0.41
46:DP:49:ARG:HG2	46:DP:50:ARG:N	2.35	0.41
47:DQ:137:TYR:HE2	56:DZ:76:LEU:CD2	2.33	0.41
47:DQ:43:THR:HG1	47:DQ:46:GLN:HG3	1.83	0.41
48:DR:67:LEU:O	48:DR:70:LEU:O	2.38	0.41
48:DR:78:LYS:O	48:DR:82:GLU:HB3	2.21	0.41
49:DS:20:ARG:HA	49:DS:20:ARG:HD3	1.70	0.41
49:DS:68:GLN:O	49:DS:71:ARG:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:7:ILE:C	50:DT:9:LEU:H	2.24	0.41
50:DT:88:ILE:CD1	50:DT:88:ILE:N	2.84	0.41
35:DA:29:U:H4'	51:DU:7:GLY:O	2.21	0.41
52:DV:36:PRO:CG	52:DV:60:GLU:OE1	2.69	0.41
54:DX:54:VAL:HA	54:DX:78:LYS:O	2.21	0.41
56:DZ:24:LEU:C	56:DZ:24:LEU:CD1	2.87	0.41
56:DZ:70:LEU:HA	56:DZ:70:LEU:HD23	1.70	0.41
56:DZ:82:ARG:NH1	56:DZ:83:PRO:O	2.54	0.41
1:AA:1011:G:H2'	1:AA:1012:U:C5'	2.51	0.41
1:AA:1118:C:O5'	9:AI:104:ARG:HG3	2.20	0.41
1:AA:1014:A:H2	1:AA:1219:U:O2	2.03	0.41
1:AA:1226:C:OP1	19:AS:81:ARG:CZ	2.69	0.41
1:AA:1425:U:H2'	1:AA:1426:C:C6	2.56	0.41
1:AA:178:C:H2'	1:AA:179:A:C8	2.52	0.41
1:AA:363:A:C6	12:AL:31:PRO:HD2	2.56	0.41
1:AA:665:A:H2'	1:AA:732:C:O2	2.21	0.41
1:AA:751:U:H2'	1:AA:752:G:H5'	2.02	0.41
1:AA:751:U:H2'	1:AA:752:G:O4'	2.21	0.41
1:AA:757:U:O2'	1:AA:879:C:H1'	2.21	0.41
1:AA:764:C:C2	1:AA:765:G:C8	3.09	0.41
1:AA:963:G:N2	10:AJ:55:LYS:CE	2.84	0.41
2:AB:23:ARG:HG3	2:AB:23:ARG:HH11	1.86	0.41
3:AC:181:ASN:HD21	3:AC:204:LEU:HB2	1.83	0.41
3:AC:37:GLN:O	3:AC:39:ILE:N	2.54	0.41
4:AD:129:ASN:HB2	4:AD:131:ARG:HH22	1.86	0.41
5:AE:51:VAL:O	5:AE:52:PRO:C	2.59	0.41
7:AG:104:LEU:O	7:AG:105:VAL:C	2.59	0.41
1:AA:876:G:O2'	8:AH:7:ALA:HB1	2.21	0.41
11:AK:124:LYS:HD2	11:AK:125:PHE:CE1	2.55	0.41
11:AK:67:ASP:OD1	11:AK:71:LYS:HE3	2.21	0.41
12:AL:113:ARG:NH1	12:AL:120:TYR:CE2	2.88	0.41
12:AL:126:LYS:HA	12:AL:126:LYS:HD2	1.78	0.41
12:AL:46:LYS:O	12:AL:47:LYS:C	2.59	0.41
13:AM:14:ARG:HG2	13:AM:14:ARG:H	1.68	0.41
15:AO:60:VAL:O	15:AO:63:ARG:N	2.54	0.41
15:AO:8:LYS:HG2	15:AO:12:ILE:HD11	2.03	0.41
18:AR:40:LEU:HA	18:AR:43:PHE:CD1	2.56	0.41
18:AR:42:ARG:O	18:AR:44:LEU:N	2.47	0.41
18:AR:52:PRO:C	18:AR:56:THR:HG23	2.38	0.41
26:B0:26:TYR:C	26:B0:67:VAL:HG11	2.41	0.41
26:B0:82:ARG:HA	26:B0:83:PRO:HD2	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:B1:75:GLU:HB2	27:B1:76:ARG:NH2	2.36	0.41
35:BA:585:G:C4	35:BA:1251:C:N4	2.89	0.41
35:BA:1797:C:H2'	35:BA:1798:U:C6	2.56	0.41
35:BA:1850:G:C5	35:BA:1851:U:C5	3.09	0.41
35:BA:1912:A:C2	35:BA:1919:A:C6	3.09	0.41
35:BA:1955:U:O2'	35:BA:1956:U:H5'	2.21	0.41
35:BA:2687:U:C2'	35:BA:2688:U:H5'	2.50	0.41
35:BA:2840:C:O2'	35:BA:2841:C:H5'	2.20	0.41
35:BA:2892:A:H2'	35:BA:2893:G:O4'	2.21	0.41
35:BA:470:A:H2'	35:BA:471:A:O4'	2.21	0.41
35:BA:670:A:H4'	35:BA:671:C:O5'	2.20	0.41
35:BA:78:A:N1	35:BA:109:G:C6	2.89	0.41
35:BA:863:A:H2	35:BA:914:C:N4	2.19	0.41
35:BA:920:G:O2'	35:BA:921:G:H5'	2.21	0.41
35:BA:945:A:C4	35:BA:2448:A:C2	3.09	0.41
36:BB:75:G:C8	36:BB:76:G:N7	2.89	0.41
36:BB:83:G:C2'	36:BB:84:C:H5'	2.51	0.41
37:BC:44:HIS:HD2	37:BC:175:VAL:HA	1.85	0.41
37:BC:44:HIS:O	37:BC:210:ARG:HA	2.21	0.41
38:BD:109:ASP:N	38:BD:195:ALA:O	2.30	0.41
39:BE:53:PRO:O	39:BE:54:GLN:C	2.59	0.41
39:BE:71:GLY:O	39:BE:72:VAL:HB	2.20	0.41
39:BE:84:PHE:CZ	39:BE:86:PRO:HB3	2.55	0.41
41:BG:150:ASP:O	41:BG:151:ALA:CB	2.69	0.41
41:BG:39:ILE:CD1	41:BG:155:MET:SD	3.09	0.41
41:BG:63:ILE:HG22	41:BG:143:GLU:HB2	2.03	0.41
42:BH:153:LYS:CD	42:BH:153:LYS:N	2.84	0.41
42:BH:20:ALA:HB1	42:BH:21:PRO:HD2	2.01	0.41
44:BN:107:LEU:HD12	44:BN:107:LEU:C	2.42	0.41
44:BN:63:THR:HG23	44:BN:64:GLY:N	2.36	0.41
45:BO:104:ARG:O	45:BO:106:LEU:N	2.54	0.41
45:BO:6:THR:O	45:BO:20:MET:HA	2.21	0.41
46:BP:74:GLU:CD	46:BP:75:ILE:HD12	2.41	0.41
48:BR:117:VAL:O	48:BR:118:GLU:CB	2.69	0.41
48:BR:59:ASP:O	48:BR:62:ALA:N	2.51	0.41
48:BR:81:ASP:N	48:BR:81:ASP:OD2	2.54	0.41
49:BS:29:PHE:CD1	49:BS:30:ARG:N	2.89	0.41
49:BS:28:VAL:O	49:BS:89:ARG:HG2	2.21	0.41
50:BT:74:ARG:HG2	50:BT:74:ARG:NH1	2.34	0.41
52:BV:61:VAL:HG21	52:BV:100:ARG:H	1.82	0.41
52:BV:40:LEU:HA	52:BV:49:THR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:BX:62:LYS:HB3	54:BX:69:TYR:H	1.86	0.41
54:BX:88:LYS:CD	54:BX:88:LYS:N	2.83	0.41
55:BY:15:VAL:O	55:BY:16:ALA:HB2	2.21	0.41
56:BZ:109:ALA:HB3	56:BZ:145:GLU:OE1	2.21	0.41
56:BZ:37:VAL:HG23	56:BZ:37:VAL:O	2.21	0.41
1:CA:1287:A:C5	1:CA:1288:A:N6	2.89	0.41
1:CA:128:G:C2'	1:CA:129:U:H5'	2.50	0.41
1:CA:1423:G:C2	1:CA:1424:C:C5	3.09	0.41
1:CA:227:G:C2	1:CA:228:A:C4	3.09	0.41
1:CA:355:C:C2'	1:CA:356:A:H5'	2.50	0.41
1:CA:376:G:H5''	16:CP:5:ARG:HB2	2.02	0.41
1:CA:939:G:C6	1:CA:940:C:N4	2.89	0.41
2:CB:113:HIS:C	2:CB:115:LEU:N	2.74	0.41
3:CC:60:ALA:O	3:CC:61:ALA:HB2	2.21	0.41
3:CC:59:ARG:HA	3:CC:63:ASN:O	2.21	0.41
4:CD:100:ARG:HG2	4:CD:100:ARG:HH11	1.85	0.41
4:CD:157:LEU:O	4:CD:161:ASN:OD1	2.38	0.41
4:CD:201:GLN:NE2	4:CD:204:ILE:HD12	2.35	0.41
6:CF:21:LEU:CA	6:CF:24:GLU:HG2	2.49	0.41
7:CG:24:THR:O	7:CG:28:ASN:ND2	2.54	0.41
8:CH:102:ARG:N	8:CH:102:ARG:HE	2.19	0.41
8:CH:109:ILE:HG23	8:CH:137:VAL:HG23	2.03	0.41
8:CH:63:LEU:HD22	8:CH:63:LEU:N	2.36	0.41
8:CH:63:LEU:CB	8:CH:65:TYR:CE1	3.04	0.41
9:CI:113:LYS:HB2	9:CI:116:LYS:CG	2.48	0.41
9:CI:88:TYR:O	9:CI:89:ASN:CB	2.69	0.41
12:CL:117:ARG:C	12:CL:119:LYS:O	2.59	0.41
12:CL:28:LYS:C	12:CL:30:ALA:H	2.24	0.41
12:CL:55:VAL:CA	12:CL:70:ILE:HD13	2.51	0.41
13:CM:92:HIS:ND1	13:CM:98:VAL:HG21	2.36	0.41
15:CO:66:LEU:O	15:CO:69:TYR:N	2.53	0.41
1:CA:1320:C:H42	19:CS:36:ARG:HD3	1.86	0.41
23:CW:31:G:N2	23:CW:42:C:C2	2.89	0.41
25:CY:152:ASP:OD1	25:CY:153:GLU:N	2.53	0.41
25:CY:2:THR:C	25:CY:4:LYS:N	2.71	0.41
25:CY:80:GLU:C	25:CY:82:ALA:N	2.74	0.41
26:D0:72:ARG:HB2	26:D0:76:GLY:O	2.21	0.41
28:D2:32:LEU:CD2	28:D2:44:LEU:HD21	2.51	0.41
35:DA:1225:G:O2'	35:DA:1226:A:H5'	2.21	0.41
35:DA:1375:C:H2'	35:DA:1376:C:H6	1.86	0.41
35:DA:1398:C:OP1	54:DX:53:LYS:NZ	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1484:G:H1	35:DA:1506:C:H42	1.69	0.41
35:DA:1761:C:H3'	35:DA:1762:A:H8	1.86	0.41
35:DA:1771:C:C1'	35:DA:1786:A:H8	2.33	0.41
35:DA:1992:G:O2'	35:DA:1993:U:P	2.79	0.41
35:DA:19:C:H2'	35:DA:20:C:C6	2.56	0.41
35:DA:202:U:N3	35:DA:203:C:C4	2.89	0.41
35:DA:2193:G:C2	35:DA:2194:G:C8	3.09	0.41
35:DA:2228:G:H2'	35:DA:2229:C:H6	1.86	0.41
35:DA:2412:A:C6	35:DA:2413:G:C4	3.09	0.41
35:DA:271(S):G:H2'	35:DA:271(T):C:O4'	2.21	0.41
35:DA:2721:A:H2'	35:DA:2722:G:H8	1.86	0.41
35:DA:2723:C:H2'	35:DA:2724:C:C5'	2.50	0.41
35:DA:2809:A:C2'	35:DA:2810:A:H5'	2.51	0.41
35:DA:38:A:C6	35:DA:39:C:C4	3.09	0.41
35:DA:709:U:C2	35:DA:723:G:N2	2.88	0.41
35:DA:910:A:C6	35:DA:911:A:C6	3.09	0.41
36:DB:32:C:H2'	36:DB:33:G:O4'	2.21	0.41
39:DE:181:LEU:N	39:DE:181:LEU:CD2	2.75	0.41
39:DE:9:VAL:HG13	39:DE:25:VAL:HG12	2.02	0.41
40:DF:20:LEU:O	40:DF:21:ALA:O	2.39	0.41
40:DF:82:ILE:C	40:DF:84:VAL:H	2.23	0.41
41:DG:104:GLU:C	41:DG:106:LEU:H	2.23	0.41
41:DG:119:GLY:O	41:DG:120:LEU:C	2.58	0.41
41:DG:12:TYR:CA	41:DG:16:ARG:HG3	2.51	0.41
42:DH:117:PRO:HA	42:DH:118:PRO:HD2	1.99	0.41
42:DH:162:ILE:C	42:DH:162:ILE:CD1	2.83	0.41
42:DH:19:VAL:HG12	42:DH:20:ALA:N	2.36	0.41
42:DH:87:LEU:CD1	42:DH:148:ILE:HG21	2.51	0.41
43:DI:26:ALA:HA	43:DI:30:LEU:HB2	2.02	0.41
44:DN:2:LYS:HZ1	52:DV:12:TYR:HB3	1.86	0.41
46:DP:17:LYS:C	46:DP:19:VAL:N	2.60	0.41
35:DA:662:G:P	46:DP:18:ARG:HD2	2.60	0.41
35:DA:911:A:N6	47:DQ:10:ARG:HG2	2.36	0.41
48:DR:80:PHE:C	48:DR:82:GLU:H	2.24	0.41
49:DS:106:ARG:NH1	49:DS:107:GLU:O	2.54	0.41
49:DS:13:ARG:N	49:DS:13:ARG:CD	2.69	0.41
50:DT:23:ARG:CZ	50:DT:120:ARG:HD3	2.51	0.41
45:DO:71:ARG:HH12	50:DT:74:ARG:NH2	2.19	0.41
53:DW:12:ILE:CD1	53:DW:42:ARG:NH1	2.83	0.41
56:DZ:28:MET:HG3	56:DZ:33:LEU:CD2	2.50	0.41
56:DZ:52:SER:O	56:DZ:71:VAL:HG21	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:8:TYR:H	56:DZ:62:PRO:HD3	1.85	0.41
1:AA:1054:C:O2	1:AA:1054:C:C2'	2.67	0.41
1:AA:1153:C:C2'	1:AA:1154:G:O5'	2.69	0.41
1:AA:1154:G:O2'	1:AA:1155:G:H5'	2.21	0.41
1:AA:1164:G:C6	1:AA:1173:G:C6	3.08	0.41
1:AA:1267:C:C2'	1:AA:1267:C:O2	2.69	0.41
1:AA:1414:U:H2'	1:AA:1415:G:H8	1.86	0.41
1:AA:349:A:C2'	1:AA:350:G:H5'	2.50	0.41
1:AA:389:A:H2'	1:AA:390:C:O4'	2.21	0.41
1:AA:437:U:H2'	1:AA:438:G:H5'	2.03	0.41
1:AA:445:G:H2'	1:AA:446:G:O4'	2.21	0.41
1:AA:482:A:N3	1:AA:482:A:H2'	2.35	0.41
1:AA:502:G:C2	1:AA:544:G:C2	3.08	0.41
1:AA:506:G:H2'	1:AA:507:C:H6	1.83	0.41
1:AA:763:G:H2'	1:AA:764:C:H6	1.85	0.41
1:AA:837:G:O2'	1:AA:838:G:H5'	2.20	0.41
1:AA:832:C:C4	1:AA:855:G:N1	2.89	0.41
2:AB:30:ARG:HG3	2:AB:31:TYR:CD2	2.55	0.41
2:AB:80:ILE:HD11	2:AB:215:LEU:CD1	2.51	0.41
3:AC:104:GLN:CD	3:AC:105:GLU:N	2.70	0.41
3:AC:188:LEU:CD2	3:AC:188:LEU:N	2.85	0.41
3:AC:206:GLU:HB3	3:AC:207:VAL:H	1.59	0.41
4:AD:61:LYS:HZ2	4:AD:62:GLN:NE2	2.19	0.41
5:AE:94:ALA:HB3	5:AE:117:ASP:O	2.21	0.41
7:AG:155:ARG:HH11	7:AG:155:ARG:HG3	1.85	0.41
7:AG:49:ILE:CG2	7:AG:49:ILE:O	2.67	0.41
8:AH:22:GLU:HA	8:AH:22:GLU:OE2	2.21	0.41
9:AI:39:GLY:O	9:AI:41:VAL:N	2.54	0.41
10:AJ:85:LEU:O	10:AJ:86:MET:C	2.59	0.41
12:AL:44:THR:HA	12:AL:45:PRO:HD3	1.85	0.41
13:AM:27:LYS:O	13:AM:30:ALA:HB3	2.21	0.41
15:AO:36:ILE:HG22	15:AO:37:ASN:N	2.36	0.41
16:AP:18:ARG:HD3	16:AP:35:LYS:CD	2.44	0.41
17:AQ:29:HIS:ND1	17:AQ:31:LEU:N	2.51	0.41
1:AA:194:C:H4'	20:AT:65:LYS:HG3	2.03	0.41
25:AY:126:ARG:HA	25:AY:169:ILE:HD11	2.02	0.41
25:AY:45:TYR:HB3	25:AY:50:VAL:CG2	2.50	0.41
26:B0:56:ASP:O	26:B0:58:THR:N	2.54	0.41
27:B1:76:ARG:HE	27:B1:76:ARG:HA	1.85	0.41
28:B2:14:ARG:C	28:B2:16:LEU:N	2.74	0.41
30:B4:21:VAL:C	30:B4:23:GLU:N	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:B6:43:CYS:O	32:B6:44:ARG:O	2.38	0.41
35:BA:1106:A:H2'	35:BA:1107:G:C8	2.55	0.41
35:BA:142:A:H1'	35:BA:1408:C:O4'	2.21	0.41
35:BA:1449:A:C6	35:BA:1450:G:H1'	2.56	0.41
35:BA:1591:G:H2'	35:BA:1592:C:H5'	2.02	0.41
35:BA:1313:U:C6	35:BA:1610:A:C2	3.09	0.41
35:BA:16:G:H2'	35:BA:17:G:C8	2.56	0.41
35:BA:1682:G:C4	35:BA:1757:U:C2	3.09	0.41
35:BA:17:G:H2'	35:BA:18:C:H6	1.86	0.41
35:BA:1925:C:H2'	35:BA:1926:U:H5'	2.01	0.41
35:BA:2183:C:O2'	35:BA:2184:G:H5'	2.21	0.41
35:BA:2485:G:C2	35:BA:2486:G:C8	3.09	0.41
35:BA:2550:G:O2'	35:BA:2551:C:H5'	2.20	0.41
35:BA:2606:C:H2'	35:BA:2607:G:H5'	2.02	0.41
35:BA:2637:U:H1'	35:BA:2782:G:H22	1.86	0.41
35:BA:271(E):U:C2	35:BA:271(F):C:C5	3.09	0.41
35:BA:2746:U:H2'	35:BA:2747:G:H8	1.86	0.41
35:BA:2805:G:C2	35:BA:2807:G:C5	3.08	0.41
35:BA:638:G:H2'	35:BA:639:U:O4'	2.21	0.41
35:BA:678:C:C2	35:BA:679:C:C5	3.09	0.41
35:BA:738:G:C2'	35:BA:739:G:H5'	2.51	0.41
35:BA:827:U:C4	35:BA:2430:A:C6	3.08	0.41
35:BA:830:G:N3	35:BA:2448:A:N6	2.68	0.41
35:BA:869:G:H2'	35:BA:870:A:O4'	2.20	0.41
35:BA:900:A:N7	35:BA:901:A:N7	2.68	0.41
35:BA:987:G:H2'	35:BA:988:A:O4'	2.21	0.41
36:BB:74:U:C3'	36:BB:75:G:C5'	2.91	0.41
35:BA:2127:G:H5'	37:BC:36:LYS:HZ3	1.86	0.41
38:BD:226:MET:HE1	38:BD:230:ASP:HB3	1.98	0.41
41:BG:128:ARG:O	41:BG:129:GLY:O	2.39	0.41
41:BG:5:VAL:HB	41:BG:104:GLU:OE1	2.21	0.41
42:BH:79:VAL:O	42:BH:81:GLU:N	2.49	0.41
43:BI:133:HIS:O	43:BI:134:PRO:C	2.59	0.41
45:BO:53:LYS:O	45:BO:54:GLU:C	2.59	0.41
47:BQ:93:TYR:N	47:BQ:93:TYR:CD1	2.89	0.41
50:BT:35:LYS:HZ1	50:BT:41:ARG:HH21	1.67	0.41
51:BU:101:ARG:O	51:BU:102:GLU:C	2.59	0.41
51:BU:36:ARG:HG2	51:BU:40:PHE:CZ	2.56	0.41
52:BV:2:PHE:CB	52:BV:42:GLY:CA	2.95	0.41
53:BW:76:VAL:CG2	53:BW:103:ILE:HG22	2.51	0.41
54:BX:47:PHE:O	54:BX:49:VAL:CG2	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:13:VAL:HG12	55:BY:14:LEU:H	1.86	0.41
56:BZ:105:VAL:H	56:BZ:141:VAL:CG1	2.33	0.41
1:CA:1023:G:C2'	1:CA:1024:G:H5'	2.50	0.41
1:CA:332:G:O2'	1:CA:333:G:H5'	2.21	0.41
1:CA:348:G:C2'	1:CA:349:A:H5'	2.51	0.41
1:CA:513:C:H2'	1:CA:514:C:C6	2.56	0.41
1:CA:502:G:C6	1:CA:544:G:N1	2.89	0.41
1:CA:934:C:HO2'	1:CA:1344:C:H5	1.65	0.41
2:CB:145:LEU:CD2	2:CB:149:LEU:HD12	2.51	0.41
2:CB:204:ASN:HB3	2:CB:210:SER:HB3	2.01	0.41
2:CB:23:ARG:HG3	2:CB:23:ARG:HH11	1.86	0.41
4:CD:102:ASP:HB3	4:CD:136:PRO:HB3	2.02	0.41
4:CD:128:VAL:CG1	4:CD:129:ASN:N	2.75	0.41
4:CD:148:VAL:O	4:CD:149:ALA:C	2.60	0.41
5:CE:99:GLY:O	5:CE:117:ASP:HA	2.20	0.41
5:CE:144:THR:O	5:CE:146:ALA:N	2.54	0.41
7:CG:120:ILE:HG22	7:CG:124:LEU:CD1	2.46	0.41
7:CG:25:ALA:O	7:CG:28:ASN:HB2	2.21	0.41
8:CH:38:ILE:HD11	8:CH:118:VAL:O	2.21	0.41
8:CH:25:ASP:OD2	8:CH:60:ARG:HD3	2.21	0.41
8:CH:49:GLU:HG2	8:CH:62:TYR:CE2	2.56	0.41
9:CI:102:LEU:C	9:CI:102:LEU:HD23	2.41	0.41
1:CA:1179:A:O2'	9:CI:103:THR:HG23	2.21	0.41
9:CI:27:THR:O	9:CI:28:VAL:HG23	2.20	0.41
9:CI:28:VAL:HG22	9:CI:63:ILE:HB	2.03	0.41
9:CI:4:TYR:HB2	9:CI:19:LEU:CD1	2.48	0.41
11:CK:29:ILE:C	11:CK:29:ILE:CD1	2.86	0.41
11:CK:33:THR:CA	11:CK:40:ILE:HG12	2.50	0.41
11:CK:84:VAL:HG22	11:CK:110:ASP:HA	2.03	0.41
12:CL:90:VAL:O	12:CL:90:VAL:HG12	2.20	0.41
1:CA:911:U:OP1	12:CL:95:GLY:HA2	2.20	0.41
13:CM:100:GLY:C	13:CM:101:GLN:HG3	2.41	0.41
10:CJ:49:VAL:HG11	14:CN:41:ARG:HB2	2.01	0.41
14:CN:4:LYS:HA	14:CN:7:ILE:HG12	2.03	0.41
18:CR:42:ARG:O	18:CR:44:LEU:N	2.46	0.41
13:CM:91:ARG:NH1	19:CS:81:ARG:HH22	1.91	0.41
20:CT:43:LEU:O	20:CT:44:ALA:C	2.59	0.41
23:CW:38:A:N7	23:CW:39:A:C6	2.89	0.41
25:CY:108:GLU:N	25:CY:111:ARG:CZ	2.84	0.41
25:CY:14:MET:O	25:CY:18:LEU:CB	2.65	0.41
27:D1:13:ILE:HD11	27:D1:14:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:D1:37:ILE:HG13	27:D1:38:SER:N	2.36	0.41
27:D1:89:GLU:C	27:D1:90:ILE:HD13	2.41	0.41
32:D6:51:GLU:OE1	32:D6:51:GLU:N	2.47	0.41
33:D7:34:ARG:NE	33:D7:39:ARG:HD2	2.36	0.41
35:DA:1112:G:O2'	35:DA:1113:U:H6	2.04	0.41
35:DA:1203:G:H3'	35:DA:1204:A:H5''	2.03	0.41
35:DA:1400:G:H2'	35:DA:1401:G:H8	1.86	0.41
35:DA:1418:G:H1	35:DA:1579:A:C5'	2.30	0.41
35:DA:1662:C:O2'	35:DA:2687:U:OP1	2.39	0.41
35:DA:1706:U:O2'	35:DA:1707:G:H5'	2.21	0.41
35:DA:2124:G:O2'	37:DC:40:THR:HA	2.21	0.41
35:DA:2538:C:C2'	35:DA:2539:C:C5'	2.97	0.41
35:DA:2615:U:H2'	35:DA:2616:C:H6	1.86	0.41
35:DA:271(V):G:N3	35:DA:271(W):G:H1'	2.36	0.41
35:DA:2737:G:H1	35:DA:2767:C:H42	1.67	0.41
35:DA:2810:A:N6	35:DA:2891:G:H1'	2.36	0.41
35:DA:2892:A:H2'	35:DA:2893:G:O4'	2.21	0.41
35:DA:736:C:H2'	35:DA:737:C:C6	2.51	0.41
35:DA:99:U:H4'	35:DA:100:G:H5'	2.03	0.41
36:DB:81:G:H4'	36:DB:81:G:OP1	2.21	0.41
38:DD:133:LEU:CB	38:DD:173:VAL:HG11	2.51	0.41
38:DD:136:ILE:HG22	38:DD:137:PRO:HD2	2.03	0.41
41:DG:128:ARG:C	41:DG:129:GLY:O	2.60	0.41
41:DG:12:TYR:HA	41:DG:16:ARG:HH11	1.81	0.41
41:DG:14:GLU:C	41:DG:17:PRO:HD2	2.41	0.41
41:DG:66:GLN:NE2	41:DG:93:THR:O	2.53	0.41
41:DG:82:LEU:HD13	41:DG:87:PRO:HA	2.03	0.41
43:DI:122:GLU:O	43:DI:126:TYR:OH	2.39	0.41
43:DI:77:LEU:CD2	43:DI:101:LEU:HD13	2.50	0.41
43:DI:93:THR:N	43:DI:96:ASP:OD2	2.54	0.41
44:DN:9:VAL:CG1	44:DN:10:GLU:H	2.18	0.41
44:DN:30:ILE:HG22	44:DN:30:ILE:O	2.21	0.41
44:DN:60:ILE:O	44:DN:61:ARG:C	2.58	0.41
46:DP:113:LYS:HG2	46:DP:115:LEU:HD23	2.04	0.41
48:DR:53:HIS:C	48:DR:56:LYS:HB2	2.41	0.41
49:DS:32:LEU:HD23	49:DS:32:LEU:HA	1.88	0.41
50:DT:112:ARG:C	50:DT:112:ARG:HD3	2.42	0.41
50:DT:29:ARG:HD2	50:DT:29:ARG:HA	1.87	0.41
50:DT:51:ARG:O	50:DT:52:ILE:HG23	2.21	0.41
50:DT:52:ILE:CG1	50:DT:52:ILE:O	2.66	0.41
50:DT:64:ARG:HA	50:DT:72:VAL:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:92:GLY:O	50:DT:94:ALA:N	2.54	0.41
51:DU:18:LEU:C	51:DU:18:LEU:HD23	2.39	0.41
51:DU:36:ARG:HG2	51:DU:40:PHE:CZ	2.56	0.41
51:DU:91:ASP:O	51:DU:92:ARG:HG2	2.20	0.41
53:DW:88:ARG:HH11	53:DW:88:ARG:HG2	1.85	0.41
56:DZ:151:HIS:O	56:DZ:171:ILE:HG12	2.20	0.41
1:AA:1232:U:H5'	9:AI:126:SER:OG	2.21	0.40
1:AA:1293:G:O2'	1:AA:1294:G:O5'	2.39	0.40
1:AA:1336:C:O4'	1:AA:1337:G:N2	2.55	0.40
1:AA:1478:C:N3	1:AA:1479:C:N4	2.68	0.40
1:AA:1480:G:N2	1:AA:1481:U:HI1'	2.36	0.40
1:AA:528:C:C5	1:AA:529:G:N7	2.90	0.40
1:AA:552:U:H4'	12:AL:87:GLY:N	2.35	0.40
1:AA:631:G:H2'	1:AA:632:A:C8	2.56	0.40
2:AB:194:PRO:HB2	2:AB:195:ASP:H	1.71	0.40
4:AD:109:GLY:O	4:AD:161:ASN:HB3	2.21	0.40
4:AD:120:LEU:CD1	4:AD:120:LEU:N	2.73	0.40
4:AD:148:VAL:O	4:AD:148:VAL:HG12	2.22	0.40
4:AD:163:GLU:O	4:AD:166:LYS:N	2.38	0.40
4:AD:12:CYS:HA	4:AD:19:LEU:H	1.86	0.40
6:AF:100:ASN:CB	18:AR:28:GLU:HA	2.50	0.40
6:AF:21:LEU:C	6:AF:23:LYS:N	2.74	0.40
7:AG:18:TYR:HD1	7:AG:18:TYR:H	1.68	0.40
7:AG:27:ILE:HD11	7:AG:43:PHE:CE2	2.56	0.40
8:AH:12:ARG:O	8:AH:15:ASN:N	2.49	0.40
8:AH:25:ASP:OD2	8:AH:60:ARG:HD3	2.21	0.40
9:AI:8:GLY:HA3	9:AI:15:ALA:HB3	2.03	0.40
9:AI:27:THR:C	9:AI:28:VAL:HG23	2.42	0.40
1:AA:1253:G:OP1	10:AJ:44:VAL:HG11	2.21	0.40
12:AL:33:ARG:HB3	12:AL:85:ILE:CG2	2.51	0.40
12:AL:6:THR:HG22	12:AL:9:GLN:NE2	2.31	0.40
1:AA:981:U:H5'	14:AN:21:TYR:OH	2.22	0.40
15:AO:48:LYS:HA	15:AO:48:LYS:HD3	1.78	0.40
1:AA:393:A:P	16:AP:12:LYS:HD3	2.60	0.40
17:AQ:22:LEU:HD12	17:AQ:23:VAL:H	1.86	0.40
19:AS:5:LEU:HD12	19:AS:10:PHE:H	1.86	0.40
20:AT:25:ARG:HH11	20:AT:25:ARG:CG	2.34	0.40
20:AT:34:LYS:O	20:AT:35:THR:C	2.59	0.40
20:AT:56:MET:O	20:AT:60:GLU:CB	2.69	0.40
20:AT:78:ALA:O	20:AT:81:LYS:N	2.54	0.40
25:AY:133:ARG:NH2	35:BA:1942:C:O4'	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:B2:23:LYS:CA	54:BX:5:TYR:CE1	3.02	0.40
33:B7:29:LYS:O	33:B7:33:ARG:N	2.45	0.40
34:B8:35:GLN:HB3	34:B8:36:LYS:H	1.67	0.40
34:B8:6:THR:HG22	34:B8:62:LEU:HB2	2.03	0.40
35:BA:1022:G:C6	35:BA:1140:C:C4	3.08	0.40
35:BA:1192:G:H2'	35:BA:1193:G:H5'	2.00	0.40
35:BA:1260:G:H2'	35:BA:1261:C:H6	1.86	0.40
31:B5:11:THR:CB	35:BA:1263:U:O3'	2.68	0.40
35:BA:1437:C:O2	35:BA:1437:C:H2'	2.21	0.40
35:BA:1441:G:H2'	35:BA:1442:G:H8	1.84	0.40
35:BA:1637:A:C6	35:BA:1638:C:N4	2.89	0.40
35:BA:1721:G:C6	35:BA:1739:U:H5'	2.56	0.40
35:BA:1763:G:C4'	35:BA:1763:G:OP1	2.69	0.40
35:BA:1794:U:H2'	35:BA:1795:C:C6	2.56	0.40
35:BA:1847:A:C2'	35:BA:1847:A:N3	2.83	0.40
35:BA:1889:A:H2'	35:BA:1890:A:O4'	2.21	0.40
35:BA:1916:A:H5'	35:BA:1917:U:OP2	2.21	0.40
35:BA:2027:G:C6	35:BA:2028:U:C4	3.09	0.40
35:BA:208:C:O2'	35:BA:209:C:H5'	2.21	0.40
35:BA:2243:U:C2	35:BA:2244:U:C5	3.09	0.40
35:BA:2287:A:N6	35:BA:2344:U:H3	2.19	0.40
35:BA:2497:A:OP2	35:BA:2497:A:C8	2.74	0.40
35:BA:2583:G:H3'	35:BA:2584:U:O2	2.21	0.40
35:BA:2600:A:HO2'	35:BA:2601:C:H5'	1.79	0.40
35:BA:260:G:N3	35:BA:260:G:H2'	2.36	0.40
35:BA:2724:C:OP1	39:BE:118:LYS:HE3	2.22	0.40
35:BA:2726:U:O2'	35:BA:2727:G:H5'	2.21	0.40
35:BA:2787:C:H1'	39:BE:61:ARG:HD3	2.02	0.40
35:BA:2861:G:C2'	35:BA:2862:G:H5'	2.51	0.40
35:BA:479:A:N6	35:BA:503:A:H61	2.20	0.40
35:BA:596:G:H2'	35:BA:597:U:O4'	2.21	0.40
35:BA:769:G:H2'	35:BA:770:G:H8	1.86	0.40
35:BA:870:A:C2	35:BA:871:U:H1'	2.57	0.40
35:BA:910:A:N9	47:BQ:13:GLN:OE1	2.54	0.40
35:BA:915:C:H2'	35:BA:916:G:C8	2.56	0.40
37:BC:146:GLY:O	37:BC:148:ASN:N	2.54	0.40
38:BD:213:ARG:O	38:BD:214:TRP:C	2.59	0.40
40:BF:136:THR:O	40:BF:137:LYS:C	2.60	0.40
40:BF:171:PRO:C	40:BF:173:VAL:H	2.24	0.40
41:BG:131:TYR:O	41:BG:159:VAL:HG22	2.22	0.40
42:BH:70:THR:HB	42:BH:71:LEU:H	1.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BH:76:VAL:O	42:BH:77:LYS:C	2.59	0.40
43:BI:1:MET:HB2	43:BI:21:VAL:O	2.21	0.40
44:BN:70:LYS:HG2	44:BN:87:LEU:HD23	2.03	0.40
45:BO:104:ARG:NH2	50:BT:33:LYS:HD2	2.36	0.40
45:BO:10:VAL:HB	45:BO:12:ASP:OD2	2.21	0.40
35:BA:2566:A:C6	45:BO:28:SER:HB2	2.56	0.40
45:BO:65:THR:HA	45:BO:82:ASN:HA	2.02	0.40
45:BO:88:ASN:HB3	45:BO:92:GLU:O	2.21	0.40
46:BP:62:LEU:HD12	46:BP:62:LEU:H	1.80	0.40
51:BU:101:ARG:C	51:BU:102:GLU:HG2	2.41	0.40
51:BU:29:SER:O	51:BU:30:LYS:HG2	2.21	0.40
51:BU:92:ARG:CG	51:BU:95:LEU:H	2.22	0.40
54:BX:77:LYS:HD3	54:BX:78:LYS:N	2.36	0.40
55:BY:9:LYS:HG3	55:BY:10:GLY:N	2.36	0.40
56:BZ:109:ALA:O	56:BZ:110:GLY:C	2.59	0.40
56:BZ:58:VAL:HG22	56:BZ:68:PRO:CB	2.50	0.40
1:CA:1154:G:H2'	1:CA:1155:G:C8	2.53	0.40
1:CA:989:C:N4	1:CA:1217:C:H42	2.14	0.40
1:CA:1413:A:C6	1:CA:1488:G:N1	2.89	0.40
1:CA:1465:C:O2'	1:CA:1466:C:H5'	2.20	0.40
1:CA:1517:G:C2'	1:CA:1518:A:O5'	2.69	0.40
1:CA:1516:G:N3	1:CA:1518:A:OP2	2.54	0.40
1:CA:189(K):U:H2'	1:CA:189(L):G:H8	1.86	0.40
1:CA:327:A:C4	1:CA:329:A:C8	3.09	0.40
1:CA:35:G:C4	1:CA:550:G:N2	2.89	0.40
1:CA:416:G:C6	1:CA:417:C:C4	3.09	0.40
1:CA:57:G:C6	1:CA:58:C:N3	2.89	0.40
1:CA:60:A:H2	1:CA:107:G:N3	2.19	0.40
1:CA:854:G:OP2	1:CA:871:U:C6	2.74	0.40
1:CA:973:G:C8	1:CA:974:A:C8	3.09	0.40
2:CB:178:ARG:NH1	2:CB:178:ARG:CG	2.84	0.40
4:CD:192:GLU:C	4:CD:194:LEU:H	2.24	0.40
6:CF:19:LEU:C	6:CF:21:LEU:N	2.73	0.40
6:CF:21:LEU:O	6:CF:22:GLU:C	2.59	0.40
7:CG:130:GLY:C	7:CG:132:GLY:H	2.24	0.40
8:CH:120:THR:HG23	8:CH:123:GLU:CD	2.41	0.40
8:CH:86:ILE:HG21	8:CH:133:LEU:HD23	2.03	0.40
10:CJ:22:LYS:HZ1	10:CJ:23:ILE:HG12	1.86	0.40
10:CJ:32:ALA:N	10:CJ:78:ASN:HD21	2.18	0.40
11:CK:67:ASP:OD1	11:CK:71:LYS:HE3	2.21	0.40
15:CO:52:SER:O	15:CO:55:GLY:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CO:85:LEU:HB2	15:CO:87:ILE:CD1	2.50	0.40
18:CR:29:PHE:H	18:CR:29:PHE:HD2	1.68	0.40
18:CR:56:THR:C	18:CR:58:LEU:HD12	2.41	0.40
21:CU:6:ARG:O	21:CU:12:LYS:HD3	2.21	0.40
22:CV:29:G:H2'	22:CV:30:A:C8	2.57	0.40
25:CY:135:GLU:O	25:CY:136:ALA:C	2.58	0.40
27:D1:60:PHE:HE2	27:D1:91:LYS:HE3	1.87	0.40
28:D2:13:ALA:O	28:D2:14:ARG:NH2	2.54	0.40
28:D2:40:SER:HB2	28:D2:41:ILE:CD1	2.45	0.40
32:D6:22:ALA:HB2	32:D6:39:TYR:CE2	2.56	0.40
35:DA:1449:A:H5'	35:DA:1450:G:OP2	2.21	0.40
35:DA:1495:A:C2	35:DA:1496:A:C4	3.09	0.40
35:DA:1511:C:H2'	35:DA:1512:U:C6	2.56	0.40
35:DA:1615:C:C5	35:DA:1617:C:C6	3.10	0.40
35:DA:1747:G:H2'	35:DA:1747(A):G:H8	1.86	0.40
35:DA:1813:G:H2'	35:DA:1814:G:O4'	2.20	0.40
35:DA:1925:C:C2'	35:DA:1926:U:C5'	2.97	0.40
35:DA:1982:C:H2'	35:DA:1983:C:H6	1.85	0.40
35:DA:2011:U:C2'	35:DA:2012:G:H5'	2.51	0.40
35:DA:2041:U:H2'	35:DA:2042:A:H8	1.87	0.40
35:DA:2248:C:H2'	35:DA:2249:U:C5'	2.50	0.40
35:DA:2560:C:H2'	35:DA:2561:A:H5'	2.02	0.40
35:DA:2603:G:H4'	35:DA:2603:G:OP2	2.21	0.40
35:DA:334:C:O2'	35:DA:335:C:P	2.78	0.40
35:DA:408:G:O2'	35:DA:409:C:H5'	2.21	0.40
35:DA:483:A:H1'	55:DY:47:LYS:CG	2.25	0.40
35:DA:591:C:C2	35:DA:592:G:C8	3.09	0.40
35:DA:651:G:HO2'	35:DA:652:C:P	2.44	0.40
35:DA:937:U:H2'	35:DA:938:G:O4'	2.21	0.40
36:DB:41:U:O4	41:DG:71:THR:HA	2.21	0.40
36:DB:90:A:N7	36:DB:91:C:H1'	2.35	0.40
37:DC:44:HIS:O	37:DC:210:ARG:HA	2.22	0.40
38:DD:24:ILE:O	38:DD:25:THR:C	2.60	0.40
38:DD:36:PRO:CB	38:DD:62:TYR:O	2.69	0.40
39:DE:133:LYS:CA	39:DE:134:ILE:HD13	2.51	0.40
39:DE:4:ILE:HD12	39:DE:5:LEU:H	1.85	0.40
40:DF:119:ARG:NH1	40:DF:119:ARG:HG2	2.36	0.40
41:DG:45:GLU:OE1	41:DG:45:GLU:C	2.59	0.40
41:DG:63:ILE:HD13	41:DG:141:PHE:CZ	2.56	0.40
43:DI:96:ASP:O	43:DI:100:ALA:N	2.54	0.40
43:DI:17:GLN:CG	43:DI:18:VAL:H	2.24	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DI:54:GLN:HA	43:DI:57:ARG:CB	2.51	0.40
44:DN:7:LYS:O	44:DN:8:GLN:C	2.58	0.40
45:DO:4:PRO:O	45:DO:5:GLN:HB2	2.20	0.40
46:DP:138:LEU:O	46:DP:138:LEU:HD22	2.20	0.40
47:DQ:55:VAL:O	47:DQ:57:HIS:N	2.55	0.40
47:DQ:12:GLN:HE21	47:DQ:73:PRO:HD3	1.86	0.40
48:DR:23:ASN:O	48:DR:24:GLN:C	2.59	0.40
48:DR:34:ILE:HG22	48:DR:35:THR:H	1.86	0.40
49:DS:51:ALA:HA	49:DS:56:LEU:CD1	2.51	0.40
51:DU:61:TRP:CE2	51:DU:94:ASN:HB2	2.56	0.40
51:DU:66:ASN:O	51:DU:70:ARG:HB2	2.21	0.40
51:DU:72:HIS:CE1	51:DU:107:ALA:HA	2.56	0.40
51:DU:94:ASN:C	51:DU:96:ALA:H	2.24	0.40
52:DV:14:VAL:HG11	52:DV:98:GLU:CG	2.49	0.40
52:DV:94:LEU:HA	52:DV:94:LEU:HD23	1.84	0.40
54:DX:60:ARG:HG2	54:DX:73:ARG:N	2.36	0.40
55:DY:28:LYS:HA	55:DY:38:ILE:HG22	2.02	0.40
55:DY:31:LEU:HA	55:DY:31:LEU:HD22	1.93	0.40
55:DY:88:LYS:O	55:DY:90:LEU:HD23	2.20	0.40
56:DZ:28:MET:HB3	56:DZ:88:PHE:HB2	2.02	0.40
1:AA:1262:C:C2	1:AA:1263:C:C5	3.09	0.40
1:AA:1375:A:H4'	7:AG:29:LYS:NZ	2.37	0.40
1:AA:137:C:H2'	1:AA:138:G:H8	1.86	0.40
1:AA:259:G:H2'	1:AA:260:G:O4'	2.21	0.40
1:AA:54:C:N4	1:AA:352:C:H2'	2.32	0.40
1:AA:376:G:C2	1:AA:377:G:C5	3.09	0.40
1:AA:402:G:C5	1:AA:403:C:C5	3.09	0.40
1:AA:402:G:C6	1:AA:403:C:C5	3.10	0.40
1:AA:607:A:C4	1:AA:608:A:C8	3.09	0.40
1:AA:699:C:H2'	1:AA:700:G:H8	1.85	0.40
1:AA:666:G:C4	1:AA:741:G:N1	2.89	0.40
1:AA:805:C:O2'	1:AA:806:C:H5'	2.22	0.40
1:AA:885:G:H2'	1:AA:886:G:C8	2.55	0.40
2:AB:105:PHE:O	2:AB:107:THR:N	2.55	0.40
2:AB:11:LEU:HD11	2:AB:217:ARG:HH22	1.86	0.40
2:AB:164:VAL:O	2:AB:186:ALA:CB	2.69	0.40
2:AB:164:VAL:CG1	2:AB:165:VAL:N	2.85	0.40
1:AA:1190:G:OP2	3:AC:5:ILE:HG23	2.20	0.40
4:AD:25:ARG:C	4:AD:27:TYR:N	2.73	0.40
4:AD:58:LEU:CD1	4:AD:62:GLN:HG3	2.51	0.40
5:AE:35:GLY:HA3	5:AE:112:LEU:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AF:80:ARG:HG3	6:AF:88:VAL:HB	2.01	0.40
8:AH:25:ASP:OD1	8:AH:25:ASP:N	2.54	0.40
9:AI:27:THR:HG23	9:AI:31:GLN:O	2.20	0.40
9:AI:49:PRO:HG3	9:AI:78:LYS:HG2	2.03	0.40
9:AI:95:LYS:HZ3	9:AI:96:LEU:CB	2.30	0.40
11:AK:15:ALA:HB1	11:AK:78:GLN:HG2	2.03	0.40
12:AL:19:ARG:HD2	12:AL:19:ARG:H	1.87	0.40
12:AL:39:VAL:HG12	12:AL:40:VAL:N	2.36	0.40
12:AL:78:GLN:O	12:AL:80:HIS:N	2.55	0.40
15:AO:30:ALA:HA	15:AO:85:LEU:HD11	2.03	0.40
16:AP:51:VAL:HG12	16:AP:52:ASP:N	2.35	0.40
16:AP:60:LEU:HG	16:AP:60:LEU:H	1.59	0.40
18:AR:56:THR:C	18:AR:58:LEU:HD12	2.40	0.40
18:AR:32:ARG:NH1	18:AR:65:ILE:HG21	2.36	0.40
19:AS:42:PRO:C	19:AS:44:MET:N	2.75	0.40
25:AY:131:ASN:C	25:AY:133:ARG:N	2.73	0.40
25:AY:133:ARG:HG3	25:AY:165:THR:OG1	2.21	0.40
25:AY:36:ALA:HA	25:AY:39:LEU:HD23	2.03	0.40
27:B1:51:VAL:N	27:B1:60:PHE:O	2.48	0.40
28:B2:50:ILE:HA	28:B2:54:LYS:HD3	2.03	0.40
29:B3:51:ALA:C	29:B3:53:LEU:H	2.24	0.40
29:B3:9:VAL:HG22	29:B3:53:LEU:O	2.21	0.40
35:BA:103:A:H2'	35:BA:104:U:H6	1.86	0.40
35:BA:1223:G:N2	35:BA:1227:G:C4	2.89	0.40
35:BA:1375:C:H2'	35:BA:1376:C:H6	1.86	0.40
35:BA:1429:G:H2'	35:BA:1430:C:H6	1.86	0.40
35:BA:1560:G:H2'	35:BA:1561:G:C8	2.56	0.40
35:BA:1608:A:C5	35:BA:1611:C:C4	3.09	0.40
35:BA:1678:G:C4	35:BA:1679:U:C5	3.10	0.40
35:BA:1778:U:H5	35:BA:1784:A:C2	2.38	0.40
35:BA:1778:U:H2'	35:BA:1784:A:N6	2.36	0.40
35:BA:1882:C:H5'	35:BA:1883:G:OP2	2.21	0.40
35:BA:1986:A:H2'	35:BA:1987:G:H5''	2.03	0.40
35:BA:2235:G:H2'	35:BA:2236:C:C6	2.56	0.40
35:BA:2377:A:H2'	35:BA:2378:A:C8	2.55	0.40
35:BA:2548:G:C6	35:BA:2561:A:N1	2.89	0.40
35:BA:253:C:H2'	35:BA:254:G:H5'	2.03	0.40
35:BA:2684:U:OP2	50:BT:53:ARG:NH2	2.50	0.40
35:BA:271(J):C:C3'	35:BA:271(K):U:C5'	2.98	0.40
35:BA:2808:U:C2'	35:BA:2809:A:C5'	2.96	0.40
35:BA:2886:G:H2'	35:BA:2887:U:C6	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:379:G:N2	35:BA:396:G:C4	2.90	0.40
35:BA:13:A:N6	35:BA:525:U:C5	2.89	0.40
35:BA:817:C:O2'	35:BA:839:U:OP1	2.40	0.40
35:BA:902:C:O2'	35:BA:903:C:H5'	2.21	0.40
35:BA:926:A:C8	35:BA:926:A:H5'	2.54	0.40
35:BA:976:C:H42	35:BA:988:A:H2	1.69	0.40
38:BD:159:ALA:HB1	38:BD:198:ASN:O	2.21	0.40
38:BD:19:ALA:O	38:BD:21:PHE:CD1	2.73	0.40
39:BE:10:GLY:HA3	50:BT:8:LYS:HZ2	1.86	0.40
39:BE:72:VAL:O	39:BE:73:GLU:C	2.59	0.40
40:BF:153:SER:O	40:BF:190:GLU:HB2	2.20	0.40
40:BF:25:PRO:HB3	40:BF:119:ARG:CD	2.52	0.40
40:BF:43:LYS:HG3	40:BF:44:ARG:N	2.36	0.40
40:BF:74:ARG:H	40:BF:74:ARG:HD2	1.86	0.40
41:BG:106:LEU:C	41:BG:108:ASN:H	2.24	0.40
41:BG:91:ARG:CD	41:BG:92:VAL:N	2.85	0.40
42:BH:26:VAL:C	42:BH:32:GLU:HG3	2.41	0.40
42:BH:99:VAL:O	42:BH:100:GLY:C	2.59	0.40
43:BI:5:LEU:C	43:BI:6:LEU:HG	2.41	0.40
44:BN:39:ARG:HG3	44:BN:39:ARG:NH1	2.36	0.40
45:BO:71:ARG:NE	45:BO:105:GLU:OE2	2.50	0.40
45:BO:2:ILE:O	45:BO:33:ALA:N	2.54	0.40
45:BO:37:ASP:HB2	45:BO:62:VAL:CG2	2.50	0.40
45:BO:13:ASN:ND2	45:BO:97:ARG:N	2.69	0.40
46:BP:123:LEU:O	46:BP:124:LYS:C	2.60	0.40
46:BP:27:HIS:C	46:BP:27:HIS:CD2	2.95	0.40
47:BQ:35:VAL:CG2	47:BQ:100:GLY:O	2.69	0.40
47:BQ:70:PRO:O	47:BQ:71:ASP:HB3	2.21	0.40
48:BR:55:ALA:HB2	48:BR:79:LEU:CD1	2.50	0.40
49:BS:28:VAL:HG21	49:BS:97:ARG:NH2	2.37	0.40
49:BS:28:VAL:HA	49:BS:37:ALA:HA	2.04	0.40
49:BS:90:GLY:HA2	49:BS:92:TYR:CG	2.56	0.40
50:BT:100:TYR:C	50:BT:102:ILE:H	2.23	0.40
50:BT:100:TYR:O	50:BT:101:PHE:C	2.60	0.40
50:BT:118:ARG:O	50:BT:121:ILE:N	2.49	0.40
50:BT:45:PHE:HE2	50:BT:63:VAL:HG23	1.86	0.40
35:BA:534:U:H4'	51:BU:46:ALA:H	1.87	0.40
51:BU:57:PHE:O	51:BU:58:ARG:C	2.60	0.40
52:BV:18:LEU:N	52:BV:18:LEU:HD13	2.36	0.40
52:BV:33:VAL:HA	52:BV:62:LEU:O	2.22	0.40
54:BX:77:LYS:HE3	54:BX:77:LYS:CA	2.49	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:BY:47:LYS:HD2	55:BY:47:LYS:N	2.37	0.40
55:BY:55:TYR:O	55:BY:56:PRO:O	2.40	0.40
1:CA:1074:G:O4'	2:CB:104:ASN:HB2	2.21	0.40
1:CA:1190:G:H5''	3:CC:3:ASN:HD21	1.86	0.40
1:CA:1226:C:C5'	19:CS:80:TYR:HE2	2.33	0.40
1:CA:978:A:C4	1:CA:1319:A:C2	3.10	0.40
1:CA:1406:U:C5	1:CA:1407:C:C5	3.09	0.40
1:CA:145:G:C2	1:CA:178:C:N3	2.89	0.40
1:CA:1471:G:O2'	1:CA:1472:U:H5'	2.21	0.40
1:CA:1501:C:OP1	1:CA:1508:G:H4'	2.21	0.40
1:CA:527:G:H2'	1:CA:528:C:C5'	2.51	0.40
1:CA:763:G:C6	1:CA:764:C:C4	3.09	0.40
2:CB:70:PHE:CD1	2:CB:163:PHE:HB3	2.56	0.40
3:CC:108:ASN:HB3	3:CC:111:LEU:HB2	2.03	0.40
4:CD:126:ILE:CG2	4:CD:127:THR:N	2.81	0.40
4:CD:145:GLU:C	4:CD:146:ILE:HD13	2.42	0.40
1:CA:922:G:H4'	5:CE:20:GLN:HA	2.03	0.40
5:CE:90:VAL:O	5:CE:91:LEU:HD12	2.21	0.40
7:CG:74:GLU:CG	7:CG:75:VAL:N	2.84	0.40
8:CH:129:VAL:CG2	8:CH:130:GLY:N	2.82	0.40
8:CH:83:ILE:CD1	8:CH:137:VAL:HG22	2.39	0.40
9:CI:107:ARG:C	9:CI:108:VAL:HG22	2.41	0.40
11:CK:125:PHE:H	11:CK:125:PHE:HD1	1.68	0.40
1:CA:523:A:H61	12:CL:53:ARG:NH2	2.19	0.40
17:CQ:59:ILE:HA	17:CQ:59:ILE:HD13	1.85	0.40
18:CR:22:VAL:O	18:CR:22:VAL:HG12	2.21	0.40
19:CS:16:LEU:HD12	19:CS:16:LEU:H	1.83	0.40
23:CW:56:U:O2'	23:CW:57:C:C5	2.74	0.40
25:CY:118:VAL:O	25:CY:119:ARG:C	2.59	0.40
25:CY:75:ALA:O	25:CY:78:ALA:HB3	2.21	0.40
27:D1:78:LYS:O	27:D1:79:GLY:C	2.59	0.40
28:D2:58:ALA:O	28:D2:60:LEU:N	2.54	0.40
33:D7:11:LYS:HG3	33:D7:12:ARG:N	2.35	0.40
34:D8:11:LYS:CG	34:D8:11:LYS:O	2.69	0.40
35:DA:99:U:H1'	35:DA:102:G:C5	2.56	0.40
35:DA:1034:G:C6	35:DA:1122:G:C4	3.09	0.40
35:DA:109:G:C4	35:DA:110:G:C8	3.09	0.40
35:DA:1109:C:N4	35:DA:1110:G:C2	2.90	0.40
35:DA:1114:G:C3'	35:DA:1115:G:C5'	2.99	0.40
35:DA:1376:C:O2'	35:DA:1377:G:H5'	2.20	0.40
35:DA:1411:C:O2'	35:DA:1412:A:P	2.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1469:A:C2'	35:DA:1470:G:H5'	2.52	0.40
35:DA:1495:A:C2	35:DA:1496:A:C2	3.09	0.40
35:DA:1623:G:H2'	35:DA:1624:G:C8	2.53	0.40
35:DA:1665:A:C2'	35:DA:1666:G:H5'	2.51	0.40
35:DA:1889:A:H2'	35:DA:1890:A:O4'	2.21	0.40
35:DA:1948:G:O2'	35:DA:1949:G:H5'	2.21	0.40
35:DA:2011:U:O2'	35:DA:2012:G:H5'	2.21	0.40
35:DA:2035:G:H4'	35:DA:2036:C:OP2	2.22	0.40
35:DA:2454:G:H2'	35:DA:2455:G:H8	1.86	0.40
35:DA:2476:A:H2'	35:DA:2477:C:C5'	2.30	0.40
35:DA:2658:C:O2'	35:DA:2659:G:H5'	2.20	0.40
35:DA:2683:C:OP1	50:DT:55:ASN:ND2	2.54	0.40
35:DA:569:U:H2'	35:DA:570:G:O4'	2.21	0.40
35:DA:998:C:P	51:DU:93:LYS:CE	3.08	0.40
36:DB:87:G:C2'	36:DB:88:C:H5''	2.50	0.40
37:DC:59:ARG:O	37:DC:62:VAL:HG22	2.21	0.40
39:DE:47:VAL:HG12	39:DE:49:LEU:CD1	2.51	0.40
41:DG:37:VAL:HB	41:DG:94:LEU:CB	2.32	0.40
35:DA:2313:C:C4'	41:DG:40:ASN:ND2	2.84	0.40
41:DG:43:LEU:N	41:DG:43:LEU:HD13	2.29	0.40
36:DB:42:C:N1	41:DG:69:ALA:HB2	2.36	0.40
41:DG:4:ASP:CB	41:DG:8:LYS:HD3	2.51	0.40
42:DH:85:LYS:HE3	42:DH:141:VAL:O	2.21	0.40
42:DH:159:GLU:O	42:DH:160:LYS:CB	2.69	0.40
43:DI:88:ILE:HG13	43:DI:123:LEU:N	2.37	0.40
35:DA:2641:G:OP1	44:DN:75:TYR:CG	2.74	0.40
45:DO:43:VAL:HG12	45:DO:43:VAL:O	2.20	0.40
45:DO:53:LYS:O	45:DO:54:GLU:C	2.59	0.40
46:DP:146:VAL:HG13	46:DP:147:LEU:N	2.37	0.40
35:DA:811:U:OP2	46:DP:24:GLY:HA2	2.20	0.40
46:DP:71:VAL:HG22	46:DP:72:PRO:CG	2.50	0.40
47:DQ:134:ARG:C	47:DQ:136:ALA:N	2.74	0.40
48:DR:116:LEU:CD2	48:DR:117:VAL:HG12	2.51	0.40
48:DR:20:LEU:O	48:DR:21:TYR:C	2.59	0.40
48:DR:87:TYR:O	48:DR:88:ARG:HB3	2.21	0.40
52:DV:83:ARG:O	52:DV:84:LYS:HD2	2.21	0.40
53:DW:106:ILE:O	53:DW:106:ILE:HG13	2.19	0.40
1:AA:1225:A:N3	1:AA:1225:A:H2'	2.36	0.40
1:AA:1438:G:C6	1:AA:1464:G:N1	2.90	0.40
1:AA:148:G:H1	1:AA:174:C:N4	2.15	0.40
1:AA:176:C:C2	1:AA:177:C:C5	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:191:G:O2'	1:AA:192:U:H5'	2.21	0.40
1:AA:321:A:O2'	1:AA:322:C:H5'	2.20	0.40
1:AA:889:A:H5'	1:AA:891:U:C1'	2.51	0.40
2:AB:15:VAL:HG23	2:AB:209:ARG:NE	2.37	0.40
2:AB:187:LEU:HD23	2:AB:202:PRO:O	2.21	0.40
3:AC:23:TYR:HA	10:AJ:11:PHE:HE1	1.81	0.40
3:AC:24:ALA:HB3	3:AC:29:TYR:CD1	2.36	0.40
4:AD:100:ARG:NH1	4:AD:100:ARG:HG2	2.36	0.40
4:AD:104:VAL:HG21	4:AD:140:VAL:CG2	2.50	0.40
4:AD:15:GLU:C	4:AD:17:VAL:N	2.75	0.40
4:AD:8:VAL:HG12	4:AD:9:CYS:N	2.36	0.40
5:AE:30:ALA:O	5:AE:46:GLY:N	2.37	0.40
5:AE:76:ILE:HG23	5:AE:77:PRO:CD	2.52	0.40
5:AE:89:ILE:CD1	5:AE:91:LEU:HD11	2.51	0.40
7:AG:63:LYS:HA	7:AG:63:LYS:HD2	1.87	0.40
7:AG:65:ALA:HA	7:AG:128:ALA:HA	2.04	0.40
11:AK:66:LEU:C	11:AK:68:ALA:N	2.74	0.40
11:AK:71:LYS:O	11:AK:73:MET:N	2.54	0.40
11:AK:72:ALA:HB1	11:AK:77:MET:CE	2.52	0.40
12:AL:47:LYS:HD3	12:AL:48:PRO:HD3	2.02	0.40
12:AL:76:ASN:C	12:AL:77:LEU:HD23	2.42	0.40
13:AM:28:ALA:O	13:AM:32:GLU:HB2	2.20	0.40
14:AN:29:ARG:HD2	14:AN:29:ARG:HA	1.93	0.40
3:AC:9:GLY:HA2	14:AN:49:HIS:O	2.21	0.40
18:AR:29:PHE:O	18:AR:29:PHE:CD2	2.75	0.40
19:AS:36:ARG:HB2	19:AS:72:GLY:HA2	2.03	0.40
19:AS:50:ALA:HA	19:AS:58:VAL:O	2.22	0.40
20:AT:37:SER:HA	20:AT:84:LEU:HD21	2.03	0.40
20:AT:57:ARG:O	20:AT:60:GLU:HB3	2.21	0.40
20:AT:73:HIS:O	20:AT:76:ALA:CB	2.69	0.40
20:AT:75:ASN:O	20:AT:76:ALA:C	2.58	0.40
20:AT:8:ARG:CD	20:AT:8:ARG:N	2.84	0.40
25:AY:104:PRO:HA	25:AY:105:PRO:HD2	1.93	0.40
25:AY:147:LEU:HD22	25:AY:149:LEU:HD21	2.04	0.40
25:AY:140:LEU:CD1	25:AY:157:ALA:HB1	2.44	0.40
25:AY:62:ASP:HB2	25:AY:63:PRO:CD	2.52	0.40
30:B4:11:PRO:C	30:B4:13:ARG:N	2.75	0.40
35:BA:1039:G:N1	35:BA:1117:G:C2	2.89	0.40
35:BA:1142(A):A:C5	35:BA:1144:G:N7	2.89	0.40
35:BA:809:G:C4'	35:BA:1254:A:H1'	2.51	0.40
27:B1:50:ARG:NH2	35:BA:1363:C:OP1	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1681:G:C8	35:BA:1681:G:OP2	2.68	0.40
35:BA:1814:G:H3'	35:BA:1815:A:C8	2.55	0.40
35:BA:1937:A:O2'	35:BA:1938:A:C5'	2.65	0.40
35:BA:2011:U:O2'	35:BA:2012:G:H5'	2.21	0.40
35:BA:1268:A:C2	35:BA:2013:A:C4	3.10	0.40
35:BA:2071:A:H2'	35:BA:2072:G:H8	1.87	0.40
35:BA:2262:U:C3'	35:BA:2263:C:H5''	2.51	0.40
35:BA:2415:G:H4'	46:BP:67:MET:N	2.35	0.40
35:BA:2578:G:N2	35:BA:2579:C:C2	2.90	0.40
35:BA:2590:A:O2'	35:BA:2591:C:C5'	2.65	0.40
35:BA:2618:G:O2'	35:BA:2619:C:H5'	2.21	0.40
35:BA:374:A:C2	35:BA:375:C:H1'	2.56	0.40
35:BA:462:C:O2'	35:BA:463:G:H5'	2.22	0.40
35:BA:709:U:C2	35:BA:723:G:N2	2.89	0.40
35:BA:829:A:N7	35:BA:2247:A:O2'	2.51	0.40
35:BA:893:C:C5	35:BA:894:C:C2	3.08	0.40
35:BA:94(A):G:H2'	35:BA:95:G:O4'	2.21	0.40
37:BC:196:LEU:C	37:BC:198:ALA:N	2.74	0.40
37:BC:36:LYS:HA	37:BC:36:LYS:HD3	1.94	0.40
38:BD:73:VAL:HG13	38:BD:119:ALA:O	2.22	0.40
38:BD:14:ARG:NH1	38:BD:14:ARG:CG	2.82	0.40
35:BA:1819:A:H5''	38:BD:158:ALA:HB3	2.03	0.40
38:BD:178:PRO:C	38:BD:180:GLY:N	2.74	0.40
38:BD:185:VAL:HG12	38:BD:189:CYS:SG	2.61	0.40
38:BD:211:ARG:O	38:BD:213:ARG:N	2.54	0.40
38:BD:20:ASP:OD1	38:BD:21:PHE:N	2.55	0.40
38:BD:92:ILE:O	38:BD:92:ILE:HD12	2.22	0.40
39:BE:52:LEU:O	39:BE:53:PRO:O	2.40	0.40
39:BE:50:GLY:HA3	39:BE:74:PRO:HG3	2.02	0.40
40:BF:119:ARG:HG2	40:BF:119:ARG:NH1	2.36	0.40
40:BF:32:LEU:C	40:BF:32:LEU:CD2	2.84	0.40
41:BG:100:TRP:O	41:BG:103:LEU:HB2	2.21	0.40
41:BG:142:PRO:O	41:BG:144:ILE:N	2.54	0.40
36:BB:45:A:C1'	41:BG:95:ARG:NH2	2.83	0.40
44:BN:62:VAL:HG22	44:BN:66:LYS:CB	2.51	0.40
44:BN:85:ILE:HA	44:BN:85:ILE:HD13	1.90	0.40
44:BN:91:LEU:O	44:BN:95:PRO:HD3	2.22	0.40
45:BO:86:ILE:H	45:BO:86:ILE:HD12	1.79	0.40
46:BP:35:HIS:HD2	46:BP:35:HIS:O	2.04	0.40
47:BQ:74:TYR:O	47:BQ:89:ASN:N	2.48	0.40
48:BR:18:LEU:CD1	48:BR:19:ALA:N	2.78	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BR:4:LEU:C	48:BR:6:SER:N	2.74	0.40
50:BT:29:ARG:HG2	50:BT:86:ILE:N	2.36	0.40
51:BU:8:VAL:O	51:BU:10:ARG:N	2.55	0.40
51:BU:91:ASP:OD2	51:BU:96:ALA:CB	2.70	0.40
52:BV:15:GLU:CB	52:BV:16:PRO:HD2	2.41	0.40
53:BW:20:VAL:HG23	53:BW:21:VAL:H	1.85	0.40
56:BZ:128:VAL:CG2	56:BZ:132:ASN:HB2	2.49	0.40
56:BZ:157:LEU:HD22	56:BZ:161:VAL:HB	2.03	0.40
1:CA:1105:A:C2	1:CA:1106:G:N7	2.89	0.40
1:CA:1418:A:N3	35:DA:1959:G:H1'	2.36	0.40
1:CA:129(A):G:H22	1:CA:189(E):U:H1'	1.86	0.40
1:CA:191:G:O2'	1:CA:192:U:H5'	2.21	0.40
1:CA:250:A:H1'	1:CA:252:U:C4	2.56	0.40
1:CA:665:A:H2'	1:CA:732:C:O2	2.22	0.40
1:CA:779:C:HO2'	1:CA:780:A:H5'	1.84	0.40
1:CA:889:A:H5'	1:CA:891:U:C1'	2.52	0.40
1:CA:244:U:C6	1:CA:894:G:N2	2.89	0.40
1:CA:992:U:H3	1:CA:1044:A:N6	2.20	0.40
2:CB:45:GLN:O	2:CB:48:MET:N	2.43	0.40
1:CA:1111:A:C2	3:CC:177:THR:HG23	2.56	0.40
1:CA:1189:C:C5'	3:CC:5:ILE:HG21	2.25	0.40
4:CD:4:TYR:CE2	4:CD:7:PRO:O	2.74	0.40
1:CA:1396:A:H2	5:CE:19:MET:HB2	1.87	0.40
6:CF:62:TRP:HB2	18:CR:35:ARG:NH1	2.31	0.40
7:CG:65:ALA:HA	7:CG:128:ALA:HA	2.03	0.40
7:CG:15:ASP:O	7:CG:19:GLY:HA2	2.21	0.40
1:CA:1291:G:H4'	9:CI:38:GLN:O	2.21	0.40
10:CJ:32:ALA:O	10:CJ:33:GLN:O	2.39	0.40
14:CN:3:ARG:CB	14:CN:3:ARG:HH11	2.35	0.40
15:CO:56:LEU:O	15:CO:59:MET:N	2.54	0.40
16:CP:40:ASP:H	16:CP:48:TRP:HB2	1.87	0.40
1:CA:376:G:H5''	16:CP:5:ARG:CG	2.51	0.40
19:CS:5:LEU:H	19:CS:6:LYS:NZ	2.19	0.40
20:CT:63:ILE:O	20:CT:64:ASP:C	2.59	0.40
25:CY:121:TYR:O	25:CY:122:ALA:C	2.59	0.40
25:CY:6:LEU:O	25:CY:9:GLU:HB2	2.20	0.40
27:D1:23:LYS:HZ2	27:D1:37:ILE:HD11	1.86	0.40
28:D2:48:HIS:CD2	35:DA:75:G:O2'	2.74	0.40
32:D6:35:GLU:HG3	32:D6:35:GLU:O	2.21	0.40
35:DA:1019:U:O2'	35:DA:1021:A:C2	2.72	0.40
35:DA:1108:U:H2'	35:DA:1109:C:C5'	2.48	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:1254:A:H5'	35:DA:1255:U:C5'	2.51	0.40
35:DA:1480:G:N3	35:DA:1480:G:H2'	2.36	0.40
35:DA:1496:A:H1'	35:DA:1578:U:H1'	2.03	0.40
35:DA:1992:G:O2'	35:DA:1993:U:O5'	2.39	0.40
35:DA:215:G:O4'	35:DA:216:A:H4'	2.22	0.40
35:DA:2219:G:H2'	35:DA:2220:G:H5'	2.02	0.40
35:DA:2669:G:N3	35:DA:2670:A:C8	2.89	0.40
35:DA:357:A:H8	35:DA:357:A:O5'	2.05	0.40
35:DA:374:A:H2'	35:DA:375:C:C5'	2.51	0.40
35:DA:54:G:C6	35:DA:55:G:C5	3.10	0.40
35:DA:58:G:H1	35:DA:69:C:N4	2.19	0.40
35:DA:708:C:O2	35:DA:708:C:H2'	2.20	0.40
35:DA:747:U:O2	35:DA:2014:A:H1'	2.20	0.40
35:DA:768:G:H2'	35:DA:769:G:C8	2.53	0.40
36:DB:10:C:C4	36:DB:11:C:C5	3.09	0.40
36:DB:111:G:C6	36:DB:112:U:C4	3.10	0.40
39:DE:52:LEU:HD23	39:DE:75:VAL:HG22	1.95	0.40
41:DG:160:VAL:O	41:DG:161:THR:OG1	2.34	0.40
42:DH:162:ILE:HD12	42:DH:163:TYR:N	2.35	0.40
42:DH:26:VAL:C	42:DH:32:GLU:HG3	2.40	0.40
42:DH:87:LEU:HA	42:DH:164:TYR:O	2.22	0.40
43:DI:100:ALA:O	43:DI:104:GLN:CD	2.59	0.40
44:DN:17:ASP:CG	44:DN:56:ASN:HB3	2.42	0.40
44:DN:66:LYS:HE3	44:DN:66:LYS:CA	2.51	0.40
45:DO:9:GLU:O	45:DO:83:ALA:HB1	2.21	0.40
46:DP:105:LEU:O	46:DP:107:LYS:HD2	2.20	0.40
46:DP:88:LEU:C	46:DP:90:ARG:N	2.73	0.40
46:DP:96:THR:HG22	46:DP:126:VAL:H	1.86	0.40
49:DS:94:TYR:O	49:DS:95:HIS:ND1	2.49	0.40
50:DT:109:GLU:C	50:DT:112:ARG:HG3	2.41	0.40
50:DT:32:TYR:HD2	50:DT:81:PRO:HB2	1.86	0.40
35:DA:997:G:OP1	51:DU:93:LYS:HE3	2.22	0.40
52:DV:71:LEU:CD1	52:DV:72:VAL:H	2.19	0.40
53:DW:13:SER:HA	53:DW:14:PRO:HD3	1.77	0.40
53:DW:14:PRO:O	53:DW:16:LYS:N	2.55	0.40
53:DW:27:LYS:O	53:DW:71:VAL:HG23	2.22	0.40
53:DW:50:VAL:HG13	53:DW:51:LEU:N	2.37	0.40
54:DX:77:LYS:HD3	54:DX:78:LYS:H	1.86	0.40
55:DY:15:VAL:HG12	55:DY:16:ALA:N	2.30	0.40
55:DY:87:LYS:O	55:DY:88:LYS:HB2	2.21	0.40
55:DY:87:LYS:HG2	55:DY:88:LYS:N	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:DZ:109:ALA:C	56:DZ:110:GLY:O	2.54	0.40
56:DZ:69:THR:O	56:DZ:70:LEU:HG	2.22	0.40
56:DZ:71:VAL:HG12	56:DZ:86:VAL:CG1	2.52	0.40
1:AA:1009:G:C2'	1:AA:1010:G:H5'	2.52	0.40
1:AA:1305:G:N2	1:AA:1331:G:O2'	2.40	0.40
1:AA:1476:G:C6	1:AA:1477:C:N4	2.90	0.40
1:AA:1506:U:O2'	1:AA:1507:A:H5'	2.22	0.40
1:AA:180:U:H2'	1:AA:181:G:C5'	2.51	0.40
1:AA:420:U:C5	1:AA:422:C:N3	2.90	0.40
1:AA:539:A:H2'	1:AA:540:G:C8	2.54	0.40
1:AA:597:G:C6	1:AA:644:G:C6	3.09	0.40
1:AA:783:C:C4	1:AA:784:C:H5	2.39	0.40
1:AA:952:U:H2'	1:AA:953:G:C8	2.57	0.40
1:AA:96:U:O2'	1:AA:97:G:P	2.80	0.40
1:AA:991:U:OP2	1:AA:991:U:H6	2.04	0.40
2:AB:145:LEU:CD2	2:AB:149:LEU:HD12	2.50	0.40
2:AB:16:HIS:CA	2:AB:210:SER:HB2	2.50	0.40
3:AC:206:GLU:O	3:AC:207:VAL:C	2.59	0.40
3:AC:9:GLY:CA	14:AN:49:HIS:O	2.69	0.40
4:AD:63:LYS:N	4:AD:66:ARG:NH1	2.69	0.40
5:AE:101:ILE:HD11	5:AE:119:LEU:HD23	2.03	0.40
5:AE:136:MET:C	5:AE:138:ALA:N	2.73	0.40
7:AG:107:ALA:HB1	7:AG:134:ALA:HB2	2.03	0.40
1:AA:1240:U:N3	7:AG:30:ILE:HG22	2.25	0.40
8:AH:35:ILE:O	8:AH:36:LEU:C	2.60	0.40
9:AI:113:LYS:HB2	9:AI:116:LYS:CG	2.47	0.40
9:AI:3:GLN:O	9:AI:88:TYR:CE1	2.75	0.40
11:AK:123:LYS:HE2	11:AK:123:LYS:HB3	1.83	0.40
12:AL:55:VAL:C	12:AL:70:ILE:HD11	2.41	0.40
1:AA:1226:C:OP2	13:AM:103:THR:HG21	2.21	0.40
13:AM:83:ASP:CG	13:AM:84:ILE:H	2.23	0.40
15:AO:28:GLN:O	15:AO:29:VAL:C	2.60	0.40
19:AS:22:LEU:C	19:AS:24:ALA:N	2.74	0.40
25:AY:153:GLU:OE2	26:B0:8:ALA:HA	2.22	0.40
26:B0:17:GLN:HG2	35:BA:2261:C:OP1	2.21	0.40
26:B0:56:ASP:OD2	26:B0:58:THR:OG1	2.37	0.40
27:B1:34:THR:O	27:B1:35:THR:CG2	2.69	0.40
28:B2:29:LYS:HA	28:B2:32:LEU:CG	2.51	0.40
28:B2:30:ARG:O	28:B2:34:GLU:HG2	2.22	0.40
29:B3:35:ARG:NH1	29:B3:35:ARG:HG3	2.34	0.40
29:B3:41:PRO:N	29:B3:44:ARG:HD2	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B7:18:PHE:O	33:B7:20:ALA:N	2.55	0.40
35:BA:1354:A:C2'	35:BA:1355:G:H5'	2.52	0.40
35:BA:1392:A:C5	35:BA:1393:A:C6	3.10	0.40
35:BA:1411:C:O2'	35:BA:1412:A:P	2.79	0.40
35:BA:1545:A:N7	35:BA:1546:C:C2	2.90	0.40
35:BA:1615:C:C5	35:BA:1617:C:C5	3.09	0.40
35:BA:1622:G:C2	35:BA:1623:G:C8	3.09	0.40
35:BA:1742:G:N7	35:BA:1743:C:C2	2.88	0.40
35:BA:1970:A:H1'	35:BA:1972:A:C8	2.56	0.40
35:BA:1992:G:O2'	35:BA:1993:U:O5'	2.35	0.40
35:BA:2029:G:C4	35:BA:2031:A:OP2	2.73	0.40
35:BA:2220:G:H2'	35:BA:2221:G:C8	2.40	0.40
35:BA:2408:U:O2'	35:BA:2409:G:H5'	2.22	0.40
35:BA:2410:G:H2'	35:BA:2411:A:C8	2.57	0.40
35:BA:2627:G:N2	35:BA:2781:A:H2	2.18	0.40
35:BA:2687:U:C4	35:BA:2688:U:C5	3.10	0.40
35:BA:2729:G:C5	35:BA:2730:C:C4	3.09	0.40
35:BA:26:G:N1	35:BA:27:G:N2	2.69	0.40
35:BA:283:A:O2'	35:BA:284:U:OP1	2.26	0.40
35:BA:325:G:H2'	35:BA:326:G:C8	2.55	0.40
35:BA:738:G:C2	35:BA:759:G:C5	3.10	0.40
35:BA:856:C:H4'	35:BA:857:C:OP1	2.21	0.40
36:BB:9:G:C2	36:BB:113:G:C5	3.10	0.40
37:BC:49:ILE:HB	37:BC:50:ASP:H	1.49	0.40
38:BD:145:VAL:O	38:BD:153:ALA:HB1	2.20	0.40
38:BD:138:VAL:CG2	38:BD:166:GLN:O	2.69	0.40
38:BD:25:THR:HG21	38:BD:81:ALA:CB	2.28	0.40
39:BE:35:GLN:HE21	39:BE:37:ARG:NE	2.20	0.40
40:BF:114:VAL:CG2	40:BF:115:ALA:N	2.71	0.40
40:BF:158:THR:HG21	40:BF:163:VAL:CG1	2.52	0.40
40:BF:41:LEU:HA	40:BF:44:ARG:CD	2.50	0.40
41:BG:138:GLN:OE1	41:BG:153:ARG:N	2.55	0.40
41:BG:47:LYS:N	41:BG:51:ARG:HG3	2.36	0.40
42:BH:89:ILE:HG12	42:BH:90:LYS:H	1.87	0.40
45:BO:55:GLY:O	45:BO:56:ASP:OD2	2.39	0.40
45:BO:8:LEU:N	45:BO:8:LEU:HD22	2.36	0.40
46:BP:123:LEU:CD2	46:BP:123:LEU:N	2.84	0.40
47:BQ:12:GLN:HE21	47:BQ:73:PRO:CD	2.33	0.40
48:BR:31:HIS:HB2	48:BR:34:ILE:CD1	2.51	0.40
48:BR:62:ALA:O	48:BR:63:ARG:C	2.60	0.40
49:BS:46:VAL:CG1	49:BS:47:THR:N	2.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BV:88:ARG:HH11	52:BV:88:ARG:HG3	1.84	0.40
53:BW:26:GLY:CA	53:BW:71:VAL:HB	2.51	0.40
54:BX:23:GLU:HG3	54:BX:24:GLY:N	2.36	0.40
55:BY:20:TYR:CZ	55:BY:42:VAL:HA	2.57	0.40
55:BY:8:LYS:H	55:BY:8:LYS:CD	2.10	0.40
56:BZ:125:LEU:C	56:BZ:126:VAL:CG2	2.89	0.40
56:BZ:57:ILE:O	56:BZ:69:THR:O	2.39	0.40
1:CA:1153:C:C2'	1:CA:1154:G:O5'	2.69	0.40
1:CA:1226:C:H5'	19:CS:80:TYR:HE2	1.86	0.40
1:CA:1240:U:OP1	7:CG:116:ALA:HB2	2.21	0.40
1:CA:1352:C:H2'	1:CA:1353:G:C8	2.56	0.40
1:CA:302:G:C4	1:CA:303:A:C8	3.08	0.40
1:CA:54:C:N4	1:CA:352:C:H2'	2.33	0.40
1:CA:429:U:H4'	1:CA:430:A:OP1	2.18	0.40
1:CA:524:G:O5'	1:CA:524:G:H8	2.05	0.40
1:CA:586:C:C2'	1:CA:587:G:H5'	2.51	0.40
2:CB:107:THR:HG23	2:CB:110:GLN:CD	2.41	0.40
2:CB:187:LEU:CD2	2:CB:187:LEU:C	2.85	0.40
2:CB:36:ARG:NE	2:CB:36:ARG:HA	2.36	0.40
3:CC:134:ILE:HD11	3:CC:153:VAL:CG2	2.48	0.40
7:CG:11:GLN:CG	7:CG:12:LEU:H	2.34	0.40
7:CG:155:ARG:HB2	7:CG:155:ARG:CZ	2.51	0.40
7:CG:24:THR:HA	7:CG:27:ILE:HD13	2.03	0.40
7:CG:42:ILE:HA	7:CG:45:ASP:CB	2.49	0.40
8:CH:95:VAL:HG23	8:CH:95:VAL:O	2.21	0.40
10:CJ:50:ILE:HD11	14:CN:41:ARG:HH11	1.86	0.40
10:CJ:5:ARG:HG3	10:CJ:73:ASP:CG	2.42	0.40
11:CK:109:VAL:HG22	18:CR:85:LEU:O	2.22	0.40
11:CK:99:GLN:C	11:CK:101:SER:N	2.73	0.40
1:CA:538:G:P	12:CL:115:LYS:HB2	2.61	0.40
19:CS:45:VAL:C	19:CS:47:HIS:N	2.75	0.40
25:CY:120:GLN:HB3	25:CY:121:TYR:H	1.56	0.40
25:CY:137:LEU:O	25:CY:140:LEU:HB3	2.22	0.40
25:CY:30:THR:C	25:CY:32:ARG:N	2.72	0.40
25:CY:52:LEU:CD2	25:CY:53:ASN:N	2.85	0.40
25:CY:68:VAL:HG21	25:CY:99:LEU:HD12	2.03	0.40
26:D0:56:ASP:C	26:D0:58:THR:N	2.75	0.40
27:D1:37:ILE:HD12	35:DA:2080:G:C5'	2.50	0.40
27:D1:88:LYS:O	27:D1:91:LYS:N	2.54	0.40
29:D3:15:TYR:O	29:D3:20:LYS:HE2	2.22	0.40
29:D3:14:GLY:H	29:D3:20:LYS:NZ	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:D3:23:LEU:N	29:D3:23:LEU:HD12	2.37	0.40
34:D8:25:MET:HB3	34:D8:26:LYS:H	1.51	0.40
34:D8:6:THR:HG22	34:D8:62:LEU:HB2	2.03	0.40
35:DA:1022:G:N2	35:DA:1142(A):A:C2	2.89	0.40
35:DA:1309:G:H2'	35:DA:1310:G:C5'	2.51	0.40
35:DA:1763:G:C4'	35:DA:1763:G:OP1	2.69	0.40
35:DA:49:A:N6	35:DA:177:G:C5	2.89	0.40
35:DA:1764:G:H1	35:DA:1988:C:H42	1.69	0.40
35:DA:205:G:O2'	35:DA:206:U:OP2	2.34	0.40
35:DA:2206:G:N3	35:DA:2206:G:H5''	2.37	0.40
35:DA:2352:A:H2'	35:DA:2353:G:H5'	2.03	0.40
35:DA:2063:C:O2	35:DA:2450:A:N1	2.54	0.40
35:DA:2472:G:H2'	35:DA:2529:G:N2	2.37	0.40
35:DA:265:A:H1'	35:DA:266:G:H1'	2.03	0.40
35:DA:271(E):U:C2	35:DA:271(F):C:C5	3.09	0.40
35:DA:2727:G:H2'	35:DA:2728:U:H6	1.86	0.40
35:DA:42:G:C2	35:DA:437:G:C2	3.10	0.40
35:DA:457:A:C8	35:DA:459:U:N3	2.89	0.40
35:DA:666:G:C5	35:DA:667:U:C5	3.10	0.40
35:DA:974:G:C2	35:DA:989:G:N3	2.89	0.40
36:DB:106:G:C2	36:DB:107:G:C8	3.09	0.40
36:DB:115:G:O4'	49:DS:47:THR:CB	2.69	0.40
38:DD:24:ILE:O	38:DD:24:ILE:CD1	2.70	0.40
38:DD:49:ILE:H	38:DD:49:ILE:HG12	1.67	0.40
38:DD:76:PRO:O	38:DD:98:VAL:HG23	2.21	0.40
39:DE:28:ALA:O	39:DE:180:ASN:OD1	2.40	0.40
40:DF:132:VAL:CG2	40:DF:133:ASN:N	2.73	0.40
40:DF:150:GLY:HA2	40:DF:172:TRP:CD2	2.56	0.40
40:DF:67:GLN:O	40:DF:68:LYS:CG	2.67	0.40
40:DF:65:TRP:CZ3	40:DF:75:HIS:CD2	3.09	0.40
41:DG:119:GLY:HA2	41:DG:179:PRO:HG2	2.02	0.40
36:DB:44:G:O3'	41:DG:95:ARG:HD2	2.21	0.40
42:DH:98:LEU:HD22	42:DH:125:VAL:HB	2.03	0.40
42:DH:65:HIS:C	42:DH:67:LEU:H	2.24	0.40
42:DH:88:LEU:N	42:DH:88:LEU:HD22	2.37	0.40
43:DI:133:HIS:CG	43:DI:134:PRO:HD2	2.55	0.40
44:DN:77:GLY:O	44:DN:78:TYR:CD2	2.75	0.40
44:DN:82:LEU:HD12	44:DN:83:LYS:H	1.85	0.40
44:DN:85:ILE:CG2	44:DN:90:MET:HG2	2.51	0.40
44:DN:95:PRO:O	44:DN:96:GLU:C	2.59	0.40
34:D8:25:MET:CG	46:DP:64:LYS:HB2	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DP:75:ILE:H	46:DP:75:ILE:CD1	2.04	0.40
47:DQ:47:ILE:HG23	47:DQ:104:PHE:HZ	1.87	0.40
47:DQ:54:MET:HG3	47:DQ:64:ILE:CD1	2.51	0.40
35:DA:869:G:HO2'	47:DQ:8:LYS:HD3	1.87	0.40
48:DR:115:GLU:O	48:DR:116:LEU:C	2.60	0.40
48:DR:29:LEU:HA	48:DR:29:LEU:HD12	1.79	0.40
49:DS:31:SER:HB3	49:DS:34:HIS:O	2.22	0.40
49:DS:46:VAL:CG1	49:DS:47:THR:N	2.84	0.40
50:DT:33:LYS:HB2	50:DT:41:ARG:HB3	2.03	0.40
50:DT:64:ARG:CB	50:DT:73:GLU:HB3	2.47	0.40
51:DU:101:ARG:O	51:DU:102:GLU:C	2.59	0.40
51:DU:111:GLU:O	51:DU:112:ARG:C	2.60	0.40
51:DU:111:GLU:O	51:DU:115:ALA:N	2.43	0.40
51:DU:57:PHE:C	51:DU:59:ARG:N	2.73	0.40
51:DU:8:VAL:HB	51:DU:9:VAL:H	1.63	0.40
52:DV:19:LYS:HZ2	52:DV:21:ARG:H	1.70	0.40
52:DV:46:VAL:HG12	52:DV:47:VAL:N	2.36	0.40
53:DW:76:VAL:CG2	53:DW:103:ILE:HG22	2.51	0.40
54:DX:29:TRP:CE3	54:DX:76:ARG:HB3	2.57	0.40
55:DY:80:GLY:O	55:DY:81:LYS:HB3	2.21	0.40
55:DY:87:LYS:C	55:DY:89:PHE:H	2.23	0.40
56:DZ:150:LEU:HB2	56:DZ:154:ASP:OD1	2.22	0.40
56:DZ:149:SER:CB	56:DZ:172:ALA:O	2.70	0.40
1:AA:101:A:O2'	1:AA:102:G:H5'	2.22	0.40
1:AA:105:G:C6	1:AA:106:C:N4	2.90	0.40
1:AA:1305:G:OP1	21:AU:2:GLY:HA3	2.22	0.40
1:AA:1330:U:C5	1:AA:1331:G:C5	3.09	0.40
1:AA:1420:C:C3'	1:AA:1420:C:C6	3.03	0.40
1:AA:1433:A:C8	1:AA:1468:A:N6	2.90	0.40
1:AA:1456:G:C2'	1:AA:1457:G:H5'	2.52	0.40
1:AA:1518:A:N1	1:AA:1519:A:C2	2.90	0.40
1:AA:240:C:H2'	1:AA:241:C:C6	2.56	0.40
1:AA:315:A:H5''	1:AA:317:G:OP2	2.21	0.40
1:AA:470:C:O2'	1:AA:471:G:H5'	2.21	0.40
1:AA:579:G:H2'	1:AA:580:U:H6	1.87	0.40
1:AA:666:G:H2'	1:AA:667:G:C8	2.55	0.40
1:AA:756:C:O2'	1:AA:757:U:H5'	2.22	0.40
2:AB:239:VAL:O	2:AB:240:GLN:HB3	2.20	0.40
3:AC:156:ARG:NH2	3:AC:159:GLY:O	2.55	0.40
3:AC:87:LEU:HA	3:AC:90:GLU:HG2	2.04	0.40
4:AD:80:GLU:O	4:AD:81:GLU:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AE:105:VAL:N	5:AE:106:PRO:HD2	2.36	0.40
5:AE:140:ARG:HE	5:AE:140:ARG:HB2	1.36	0.40
7:AG:11:GLN:CG	7:AG:12:LEU:N	2.85	0.40
8:AH:91:ARG:NH1	8:AH:91:ARG:CG	2.79	0.40
9:AI:93:ARG:O	9:AI:95:LYS:N	2.50	0.40
10:AJ:51:ARG:HG2	14:AN:45:ARG:NH1	2.36	0.40
11:AK:20:TYR:O	11:AK:31:THR:N	2.52	0.40
11:AK:66:LEU:O	11:AK:70:LYS:N	2.47	0.40
15:AO:37:ASN:O	15:AO:40:SER:HB3	2.21	0.40
15:AO:51:HIS:O	15:AO:52:SER:C	2.60	0.40
17:AQ:45:HIS:C	17:AQ:73:VAL:HG23	2.42	0.40
19:AS:48:THR:HG22	19:AS:61:TYR:CB	2.51	0.40
19:AS:48:THR:HA	19:AS:60:VAL:O	2.22	0.40
25:AY:29:ARG:NH1	25:AY:110:ARG:HH21	2.19	0.40
26:B0:7:LEU:HA	47:BQ:83:MET:SD	2.61	0.40
27:B1:23:LYS:HD2	27:B1:23:LYS:HA	1.80	0.40
35:BA:1011:G:C6	35:BA:1151:G:C6	3.09	0.40
35:BA:1161:C:H2'	35:BA:1162:G:H8	1.87	0.40
35:BA:1203:G:N2	35:BA:1243:G:C6	2.90	0.40
35:BA:1369:G:N2	35:BA:1370:C:C2	2.89	0.40
35:BA:1455:G:O2'	35:BA:1456:G:H5'	2.21	0.40
35:BA:1464:C:O2'	35:BA:1465:G:H5'	2.21	0.40
35:BA:1528(A):A:H3'	35:BA:1529:G:C5'	2.46	0.40
35:BA:1653:G:H4'	35:BA:1654:A:O5'	2.21	0.40
35:BA:1663:C:HO2'	35:BA:1664:A:C5'	2.35	0.40
35:BA:1813:G:H2'	35:BA:1814:G:O4'	2.22	0.40
35:BA:1822:G:O2'	35:BA:1823:G:H5'	2.22	0.40
35:BA:1863:G:H2'	35:BA:1864:U:O4'	2.21	0.40
35:BA:2025:C:H2'	35:BA:2026:C:H6	1.87	0.40
35:BA:2080:G:H2'	35:BA:2081:C:C6	2.56	0.40
35:BA:2127:G:C1'	35:BA:2128:C:H4'	2.49	0.40
35:BA:2283:C:C5	35:BA:2389:G:C4	3.10	0.40
35:BA:2691:C:C4	35:BA:2719:G:N2	2.90	0.40
35:BA:2795:G:C6	35:BA:2802:G:N2	2.89	0.40
35:BA:328:U:H3	35:BA:332:A:H62	1.70	0.40
35:BA:366:C:H5'	35:BA:370:G:H5'	2.03	0.40
35:BA:687:C:N3	35:BA:788:A:H5'	2.36	0.40
35:BA:872:A:H4'	47:BQ:66:ILE:HD11	2.03	0.40
35:BA:983:A:H2'	35:BA:984:A:C8	2.57	0.40
37:BC:59:ARG:CB	37:BC:62:VAL:HG22	2.41	0.40
39:BE:101:ARG:CD	39:BE:169:ASN:HD22	2.27	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BE:3:GLY:O	39:BE:4:ILE:CG2	2.69	0.40
40:BF:150:GLY:C	40:BF:152:GLU:N	2.75	0.40
40:BF:183:VAL:HA	40:BF:186:ILE:HD12	2.04	0.40
40:BF:84:VAL:HB	40:BF:85:GLY:H	1.49	0.40
41:BG:50:ALA:HB1	41:BG:52:ILE:HD11	2.03	0.40
42:BH:74:ASN:C	42:BH:76:VAL:N	2.72	0.40
43:BI:77:LEU:HD12	43:BI:140:LEU:HD22	2.02	0.40
46:BP:50:ARG:CZ	46:BP:51:PHE:CE2	3.05	0.40
47:BQ:141:GLN:CD	56:BZ:70:LEU:HB3	2.42	0.40
47:BQ:28:ALA:HB3	47:BQ:105:GLU:OE1	2.22	0.40
47:BQ:55:VAL:C	47:BQ:57:HIS:N	2.72	0.40
47:BQ:85:LYS:CG	47:BQ:86:GLY:N	2.76	0.40
48:BR:116:LEU:HD22	48:BR:117:VAL:HG12	2.04	0.40
50:BT:100:TYR:H	50:BT:100:TYR:HD1	1.66	0.40
50:BT:51:ARG:HG3	50:BT:51:ARG:NH1	2.35	0.40
50:BT:29:ARG:CG	50:BT:85:LYS:HA	2.51	0.40
51:BU:55:ARG:HA	51:BU:58:ARG:HG3	2.04	0.40
52:BV:25:LEU:N	52:BV:94:LEU:HD11	2.35	0.40
53:BW:111:HIS:CG	53:BW:112:GLY:H	2.39	0.40
54:BX:43:VAL:C	54:BX:45:THR:N	2.72	0.40
55:BY:80:GLY:O	55:BY:81:LYS:HB3	2.22	0.40
56:BZ:141:VAL:O	56:BZ:141:VAL:HG22	2.20	0.40
56:BZ:148:ASP:OD1	56:BZ:149:SER:N	2.40	0.40
56:BZ:85:HIS:CG	56:BZ:86:VAL:N	2.89	0.40
1:CA:1000:U:H2'	1:CA:1001:A:H8	1.86	0.40
1:CA:1349:A:H2'	1:CA:1350:A:O5'	2.22	0.40
1:CA:1370:G:O2'	1:CA:1371:G:H5'	2.22	0.40
1:CA:1464:G:N2	1:CA:1465:C:C4	2.90	0.40
1:CA:1529:G:H3'	1:CA:1530:G:C5'	2.51	0.40
1:CA:292:G:H2'	1:CA:293:G:O4'	2.22	0.40
1:CA:318:G:C2	1:CA:319:G:N7	2.89	0.40
1:CA:38:G:C4	1:CA:397:A:N1	2.90	0.40
1:CA:509:A:C2	1:CA:510:A:C2	3.10	0.40
1:CA:684:A:H2'	1:CA:685:G:C8	2.56	0.40
1:CA:797:C:H2'	1:CA:798:G:H8	1.86	0.40
2:CB:71:VAL:HG12	2:CB:72:GLY:N	2.36	0.40
3:CC:29:TYR:HE2	14:CN:37:PHE:CE1	2.39	0.40
3:CC:9:GLY:CA	14:CN:49:HIS:O	2.70	0.40
4:CD:150:GLU:HA	4:CD:153:ARG:HD3	2.02	0.40
4:CD:80:GLU:O	4:CD:83:SER:N	2.55	0.40
5:CE:53:LEU:O	5:CE:54:ALA:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CF:29:ALA:O	6:CF:30:LEU:C	2.60	0.40
7:CG:122:HIS:O	7:CG:123:GLU:C	2.60	0.40
9:CI:66:ARG:HH11	9:CI:66:ARG:HB3	1.86	0.40
9:CI:99:LEU:HD22	9:CI:99:LEU:HA	1.89	0.40
10:CJ:20:ALA:O	10:CJ:24:VAL:HG23	2.21	0.40
11:CK:66:LEU:HB2	11:CK:67:ASP:H	1.63	0.40
12:CL:113:ARG:O	12:CL:122:THR:HG21	2.21	0.40
15:CO:31:LEU:O	15:CO:34:LEU:N	2.54	0.40
16:CP:39:TYR:HB2	16:CP:49:LEU:HB2	2.03	0.40
17:CQ:10:VAL:HG12	17:CQ:53:LEU:CD1	2.52	0.40
18:CR:24:ALA:C	18:CR:26:LEU:H	2.25	0.40
18:CR:52:PRO:C	18:CR:56:THR:HG23	2.41	0.40
19:CS:48:THR:HA	19:CS:60:VAL:O	2.21	0.40
23:CW:54:G:C2	23:CW:55:5MU:C4	3.10	0.40
25:CY:143:LEU:O	25:CY:146:GLU:CB	2.69	0.40
25:CY:169:ILE:O	25:CY:170:ALA:O	2.40	0.40
28:D2:51:ARG:NH1	35:DA:72:U:H5'	2.37	0.40
32:D6:17:LYS:O	32:D6:18:ARG:HD3	2.22	0.40
33:D7:28:ARG:O	33:D7:29:LYS:C	2.60	0.40
35:DA:9:U:O2'	35:DA:10:G:O5'	2.34	0.40
35:DA:1504:C:O2'	35:DA:1505:C:H5'	2.20	0.40
35:DA:1593:G:C6	35:DA:1594:G:C6	3.10	0.40
35:DA:1681:G:HO2'	35:DA:1762:A:H2'	1.86	0.40
35:DA:1802:A:C6	35:DA:1817:G:N2	2.88	0.40
35:DA:1997:G:C2	35:DA:1998:G:N7	2.89	0.40
35:DA:2018:G:C4	35:DA:2019:A:C8	3.10	0.40
35:DA:2023:G:C5'	35:DA:2617:C:H4'	2.52	0.40
35:DA:2228:G:C5	35:DA:2229:C:C4	3.10	0.40
35:DA:2064:C:H4'	35:DA:2251:G:N2	2.36	0.40
35:DA:2290:G:H1	35:DA:2342:C:H42	1.69	0.40
35:DA:2322:A:O2'	35:DA:2323:G:H5'	2.22	0.40
35:DA:2419:U:O2'	35:DA:2420:C:H5'	2.22	0.40
35:DA:2470:G:OP1	47:DQ:56:ARG:NH2	2.55	0.40
35:DA:2516:G:C6	35:DA:2517:C:N4	2.90	0.40
35:DA:2553:G:H2'	35:DA:2554:U:C4'	2.52	0.40
35:DA:261:G:C2	35:DA:262:A:N7	2.89	0.40
35:DA:2697:G:H2'	35:DA:2698:U:O4'	2.22	0.40
35:DA:2714:G:C6	35:DA:2715:C:C4	3.09	0.40
35:DA:2851:A:H2'	35:DA:2852:G:H8	1.87	0.40
35:DA:28:A:H2'	35:DA:28:A:N3	2.37	0.40
35:DA:385:C:O2'	35:DA:388:G:N2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DA:447:A:C2	35:DA:473:G:C8	3.09	0.40
35:DA:481:G:HO2'	35:DA:482:A:P	2.44	0.40
35:DA:806:C:H6	35:DA:806:C:O5'	2.05	0.40
35:DA:962:G:C2'	35:DA:963:U:H5'	2.52	0.40
38:DD:165:ILE:N	38:DD:165:ILE:CD1	2.71	0.40
38:DD:177:LEU:HB3	38:DD:178:PRO:CD	2.51	0.40
35:DA:2810:A:O2'	39:DE:61:ARG:CZ	2.69	0.40
40:DF:175:THR:HG23	40:DF:175:THR:O	2.21	0.40
40:DF:20:LEU:HB2	40:DF:24:LEU:HD21	2.04	0.40
40:DF:41:LEU:HA	40:DF:44:ARG:CG	2.51	0.40
40:DF:57:VAL:HG21	40:DF:87:GLY:CA	2.52	0.40
40:DF:84:VAL:HB	40:DF:85:GLY:H	1.45	0.40
42:DH:83:TYR:CD1	42:DH:84:SER:N	2.90	0.40
43:DI:97:ILE:CD1	43:DI:116:LEU:HD22	2.51	0.40
43:DI:55:ALA:O	43:DI:59:ALA:CB	2.69	0.40
44:DN:137:LYS:CG	44:DN:138:LEU:N	2.83	0.40
45:DO:87:ILE:HG23	45:DO:88:ASN:O	2.21	0.40
47:DQ:58:PHE:HD1	47:DQ:58:PHE:O	2.05	0.40
48:DR:13:HIS:O	48:DR:14:SER:C	2.60	0.40
52:DV:61:VAL:CG1	52:DV:62:LEU:H	2.30	0.40
53:DW:29:LEU:HD21	53:DW:33:ARG:NH2	2.36	0.40
53:DW:9:TYR:H	53:DW:102:HIS:HD2	1.67	0.40
56:DZ:38:TYR:CD1	56:DZ:38:TYR:O	2.75	0.40
47:DQ:140:ALA:CB	56:DZ:99:TYR:H	2.33	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BA:1593:G:O2'	36:BB:54:G:OP1[1_655]	2.17	0.03

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	AB	233/256 (91%)	114 (49%)	69 (30%)	50 (22%)	0	1
2	CB	233/256 (91%)	114 (49%)	68 (29%)	51 (22%)	0	1
3	AC	205/239 (86%)	119 (58%)	50 (24%)	36 (18%)	0	2
3	CC	205/239 (86%)	121 (59%)	47 (23%)	37 (18%)	0	2
4	AD	206/209 (99%)	101 (49%)	62 (30%)	43 (21%)	0	1
4	CD	206/209 (99%)	101 (49%)	61 (30%)	44 (21%)	0	1
5	AE	149/162 (92%)	85 (57%)	44 (30%)	20 (13%)	0	4
5	CE	149/162 (92%)	84 (56%)	45 (30%)	20 (13%)	0	4
6	AF	99/101 (98%)	61 (62%)	23 (23%)	15 (15%)	0	3
6	CF	99/101 (98%)	60 (61%)	24 (24%)	15 (15%)	0	3
7	AG	153/156 (98%)	97 (63%)	31 (20%)	25 (16%)	0	2
7	CG	153/156 (98%)	97 (63%)	31 (20%)	25 (16%)	0	2
8	AH	136/138 (99%)	78 (57%)	39 (29%)	19 (14%)	0	3
8	CH	136/138 (99%)	79 (58%)	37 (27%)	20 (15%)	0	3
9	AI	121/128 (94%)	80 (66%)	27 (22%)	14 (12%)	0	5
9	CI	121/128 (94%)	79 (65%)	27 (22%)	15 (12%)	0	5
10	AJ	97/105 (92%)	64 (66%)	23 (24%)	10 (10%)	0	7
10	CJ	97/105 (92%)	64 (66%)	22 (23%)	11 (11%)	0	6
11	AK	117/129 (91%)	69 (59%)	31 (26%)	17 (14%)	0	3
11	CK	117/129 (91%)	68 (58%)	32 (27%)	17 (14%)	0	3
12	AL	123/135 (91%)	78 (63%)	28 (23%)	17 (14%)	0	3
12	CL	123/135 (91%)	77 (63%)	27 (22%)	19 (15%)	0	3
13	AM	113/126 (90%)	59 (52%)	35 (31%)	19 (17%)	0	2
13	CM	113/126 (90%)	60 (53%)	34 (30%)	19 (17%)	0	2
14	AN	58/61 (95%)	38 (66%)	12 (21%)	8 (14%)	0	3
14	CN	58/61 (95%)	38 (66%)	12 (21%)	8 (14%)	0	3
15	AO	86/89 (97%)	48 (56%)	26 (30%)	12 (14%)	0	3
15	CO	86/89 (97%)	49 (57%)	26 (30%)	11 (13%)	0	4
16	AP	82/88 (93%)	45 (55%)	24 (29%)	13 (16%)	0	2
16	CP	82/88 (93%)	45 (55%)	24 (29%)	13 (16%)	0	2
17	AQ	98/105 (93%)	66 (67%)	19 (19%)	13 (13%)	0	4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
17	CQ	98/105 (93%)	65 (66%)	20 (20%)	13 (13%)	0	4
18	AR	68/88 (77%)	38 (56%)	18 (26%)	12 (18%)	0	2
18	CR	68/88 (77%)	37 (54%)	20 (29%)	11 (16%)	0	2
19	AS	77/93 (83%)	49 (64%)	19 (25%)	9 (12%)	0	5
19	CS	77/93 (83%)	49 (64%)	19 (25%)	9 (12%)	0	5
20	AT	97/106 (92%)	46 (47%)	29 (30%)	22 (23%)	0	1
20	CT	97/106 (92%)	46 (47%)	30 (31%)	21 (22%)	0	1
21	AU	23/27 (85%)	17 (74%)	4 (17%)	2 (9%)	1	9
21	CU	23/27 (85%)	17 (74%)	4 (17%)	2 (9%)	1	9
25	AY	183/185 (99%)	115 (63%)	44 (24%)	24 (13%)	0	4
25	CY	183/185 (99%)	81 (44%)	54 (30%)	48 (26%)	0	0
26	B0	83/85 (98%)	61 (74%)	13 (16%)	9 (11%)	0	6
26	D0	83/85 (98%)	59 (71%)	15 (18%)	9 (11%)	0	6
27	B1	87/98 (89%)	34 (39%)	22 (25%)	31 (36%)	0	0
27	D1	87/98 (89%)	41 (47%)	22 (25%)	24 (28%)	0	0
28	B2	49/72 (68%)	14 (29%)	21 (43%)	14 (29%)	0	0
28	D2	49/72 (68%)	16 (33%)	20 (41%)	13 (26%)	0	0
29	B3	58/60 (97%)	34 (59%)	15 (26%)	9 (16%)	0	3
29	D3	58/60 (97%)	34 (59%)	15 (26%)	9 (16%)	0	3
30	B4	48/71 (68%)	8 (17%)	23 (48%)	17 (35%)	0	0
30	D4	48/71 (68%)	8 (17%)	22 (46%)	18 (38%)	0	0
31	B5	57/60 (95%)	28 (49%)	14 (25%)	15 (26%)	0	0
31	D5	57/60 (95%)	28 (49%)	12 (21%)	17 (30%)	0	0
32	B6	41/54 (76%)	21 (51%)	11 (27%)	9 (22%)	0	1
32	D6	41/54 (76%)	21 (51%)	11 (27%)	9 (22%)	0	1
33	B7	47/49 (96%)	29 (62%)	9 (19%)	9 (19%)	0	2
33	D7	47/49 (96%)	29 (62%)	10 (21%)	8 (17%)	0	2
34	B8	62/65 (95%)	25 (40%)	22 (36%)	15 (24%)	0	0
34	D8	62/65 (95%)	26 (42%)	21 (34%)	15 (24%)	0	0
37	BC	183/229 (80%)	79 (43%)	49 (27%)	55 (30%)	0	0
37	DC	183/229 (80%)	79 (43%)	49 (27%)	55 (30%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	BD	270/276 (98%)	158 (58%)	62 (23%)	50 (18%)	0	2
38	DD	270/276 (98%)	159 (59%)	62 (23%)	49 (18%)	0	2
39	BE	203/206 (98%)	110 (54%)	46 (23%)	47 (23%)	0	1
39	DE	203/206 (98%)	111 (55%)	44 (22%)	48 (24%)	0	1
40	BF	206/210 (98%)	109 (53%)	53 (26%)	44 (21%)	0	1
40	DF	206/210 (98%)	108 (52%)	53 (26%)	45 (22%)	0	1
41	BG	177/182 (97%)	89 (50%)	48 (27%)	40 (23%)	0	1
41	DG	177/182 (97%)	89 (50%)	41 (23%)	47 (27%)	0	0
42	BH	158/180 (88%)	78 (49%)	44 (28%)	36 (23%)	0	1
42	DH	158/180 (88%)	80 (51%)	44 (28%)	34 (22%)	0	1
43	BI	144/148 (97%)	80 (56%)	36 (25%)	28 (19%)	0	2
43	DI	144/148 (97%)	80 (56%)	37 (26%)	27 (19%)	0	2
44	BN	137/140 (98%)	64 (47%)	40 (29%)	33 (24%)	0	0
44	DN	137/140 (98%)	65 (47%)	39 (28%)	33 (24%)	0	0
45	BO	120/122 (98%)	73 (61%)	24 (20%)	23 (19%)	0	2
45	DO	120/122 (98%)	74 (62%)	24 (20%)	22 (18%)	0	2
46	BP	144/150 (96%)	77 (54%)	27 (19%)	40 (28%)	0	0
46	DP	144/150 (96%)	75 (52%)	29 (20%)	40 (28%)	0	0
47	BQ	134/141 (95%)	70 (52%)	37 (28%)	27 (20%)	0	1
47	DQ	134/141 (95%)	72 (54%)	35 (26%)	27 (20%)	0	1
48	BR	115/118 (98%)	54 (47%)	26 (23%)	35 (30%)	0	0
48	DR	115/118 (98%)	53 (46%)	29 (25%)	33 (29%)	0	0
49	BS	97/112 (87%)	34 (35%)	27 (28%)	36 (37%)	0	0
49	DS	97/112 (87%)	34 (35%)	28 (29%)	35 (36%)	0	0
50	BT	136/146 (93%)	79 (58%)	26 (19%)	31 (23%)	0	1
50	DT	136/146 (93%)	78 (57%)	29 (21%)	29 (21%)	0	1
51	BU	115/118 (98%)	50 (44%)	39 (34%)	26 (23%)	0	1
51	DU	115/118 (98%)	51 (44%)	38 (33%)	26 (23%)	0	1
52	BV	97/101 (96%)	39 (40%)	21 (22%)	37 (38%)	0	0
52	DV	97/101 (96%)	39 (40%)	22 (23%)	36 (37%)	0	0
53	BW	111/113 (98%)	62 (56%)	27 (24%)	22 (20%)	0	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
53	DW	111/113 (98%)	60 (54%)	34 (31%)	17 (15%)	0	3
54	BX	91/96 (95%)	39 (43%)	24 (26%)	28 (31%)	0	0
54	DX	91/96 (95%)	38 (42%)	25 (28%)	28 (31%)	0	0
55	BY	99/110 (90%)	36 (36%)	29 (29%)	34 (34%)	0	0
55	DY	99/110 (90%)	38 (38%)	25 (25%)	36 (36%)	0	0
56	BZ	175/206 (85%)	87 (50%)	41 (23%)	47 (27%)	0	0
56	DZ	175/206 (85%)	92 (53%)	46 (26%)	37 (21%)	0	1
All	All	11936/12888 (93%)	6356 (53%)	3078 (26%)	2502 (21%)	0	1

All (2502) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	AB	20	GLU
2	AB	22	LYS
2	AB	29	ALA
2	AB	30	ARG
2	AB	37	ASN
2	AB	75	LYS
2	AB	77	ALA
2	AB	88	ALA
2	AB	154	LEU
2	AB	165	VAL
2	AB	167	PRO
2	AB	168	THR
2	AB	191	ASP
2	AB	194	PRO
2	AB	202	PRO
2	AB	209	ARG
2	AB	229	VAL
2	AB	238	LEU
2	AB	239	VAL
3	AC	4	LYS
3	AC	12	LEU
3	AC	53	ALA
3	AC	61	ALA
3	AC	157	ILE
3	AC	172	ARG
3	AC	174	PRO
3	AC	191	THR
3	AC	207	VAL

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Mol	Chain	Res	Type
4	AD	4	TYR
4	AD	5	ILE
4	AD	6	GLY
4	AD	9	CYS
4	AD	14	ARG
4	AD	24	GLU
4	AD	28	SER
4	AD	119	GLN
4	AD	128	VAL
4	AD	149	ALA
4	AD	150	GLU
4	AD	163	GLU
4	AD	164	ALA
4	AD	177	ASP
4	AD	191	ARG
5	AE	37	ARG
5	AE	129	ILE
6	AF	13	ASN
6	AF	69	GLU
6	AF	70	ASP
6	AF	72	VAL
7	AG	31	MET
7	AG	39	ALA
7	AG	153	HIS
8	AH	17	THR
8	AH	18	ARG
8	AH	22	GLU
8	AH	29	SER
8	AH	97	VAL
8	AH	129	VAL
9	AI	29	ASN
9	AI	117	HIS
9	AI	120	ARG
10	AJ	33	GLN
10	AJ	57	LYS
10	AJ	91	PRO
11	AK	27	ASN
11	AK	105	VAL
12	AL	10	LEU
12	AL	23	LYS
12	AL	28	LYS
12	AL	47	LYS

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Mol	Chain	Res	Type
12	AL	48	PRO
12	AL	62	SER
13	AM	4	ILE
13	AM	11	ARG
13	AM	60	VAL
13	AM	68	GLY
13	AM	83	ASP
13	AM	90	LEU
13	AM	117	VAL
14	AN	22	THR
14	AN	23	ARG
15	AO	16	ALA
17	AQ	49	GLU
17	AQ	74	LEU
17	AQ	81	ARG
17	AQ	93	GLN
17	AQ	96	GLU
18	AR	20	ALA
18	AR	37	VAL
18	AR	45	SER
19	AS	10	PHE
19	AS	67	VAL
19	AS	80	TYR
19	AS	81	ARG
20	AT	28	ALA
20	AT	49	ALA
20	AT	63	ILE
20	AT	64	ASP
25	AY	29	ARG
25	AY	69	GLN
25	AY	81	LYS
25	AY	85	ASP
25	AY	107	THR
26	B0	17	GLN
26	B0	55	ARG
26	B0	84	LEU
27	B1	11	ARG
27	B1	15	ALA
27	B1	26	ARG
27	B1	46	LEU
27	B1	49	VAL
27	B1	52	ARG

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Mol	Chain	Res	Type
27	B1	53	VAL
27	B1	57	GLU
27	B1	62	VAL
27	B1	63	ALA
27	B1	69	LYS
27	B1	75	GLU
27	B1	77	ALA
27	B1	86	SER
27	B1	87	PRO
27	B1	95	LEU
28	B2	40	SER
28	B2	42	GLY
28	B2	50	ILE
28	B2	51	ARG
28	B2	52	ASP
28	B2	59	ARG
29	B3	3	ARG
29	B3	13	ILE
29	B3	31	LEU
29	B3	51	ALA
29	B3	52	HIS
29	B3	56	VAL
29	B3	57	GLU
30	B4	6	HIS
30	B4	7	PRO
30	B4	10	VAL
30	B4	13	ARG
30	B4	14	ILE
30	B4	23	GLU
30	B4	29	PRO
30	B4	35	VAL
30	B4	47	GLN
31	B5	21	SER
31	B5	22	HIS
31	B5	32	PRO
31	B5	33	CYS
31	B5	34	PRO
31	B5	35	GLU
31	B5	57	VAL
32	B6	16	CYS
32	B6	31	PRO
32	B6	44	ARG

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Mol	Chain	Res	Type
32	B6	45	LYS
33	B7	18	PHE
33	B7	46	VAL
34	B8	4	MET
34	B8	25	MET
34	B8	32	LEU
34	B8	37	SER
34	B8	43	GLN
37	BC	22	ILE
37	BC	47	LEU
37	BC	49	ILE
37	BC	54	SER
37	BC	68	LEU
37	BC	104	LEU
37	BC	126	LYS
37	BC	133	PRO
37	BC	140	PRO
37	BC	167	LYS
37	BC	173	ALA
37	BC	174	PRO
37	BC	184	LYS
37	BC	204	ALA
37	BC	222	VAL
38	BD	3	VAL
38	BD	8	PRO
38	BD	13	ARG
38	BD	22	SER
38	BD	25	THR
38	BD	28	GLU
38	BD	32	SER
38	BD	45	ASN
38	BD	58	HIS
38	BD	135	PHE
38	BD	159	ALA
38	BD	204	ILE
38	BD	216	GLY
38	BD	223	GLY
38	BD	224	ALA
38	BD	228	PRO
38	BD	241	PRO
38	BD	262	ARG
38	BD	263	ARG

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Mol	Chain	Res	Type
38	BD	266	SER
38	BD	272	ALA
39	BE	2	LYS
39	BE	4	ILE
39	BE	45	THR
39	BE	53	PRO
39	BE	57	LYS
39	BE	61	ARG
39	BE	66	HIS
39	BE	77	ILE
39	BE	82	ARG
39	BE	88	GLY
39	BE	90	THR
39	BE	93	VAL
39	BE	122	PHE
39	BE	129	HIS
39	BE	182	LEU
39	BE	186	GLY
40	BF	2	LYS
40	BF	6	VAL
40	BF	10	PRO
40	BF	21	ALA
40	BF	40	GLN
40	BF	57	VAL
40	BF	60	SER
40	BF	66	PRO
40	BF	72	ARG
40	BF	73	ALA
40	BF	85	GLY
40	BF	89	VAL
40	BF	184	TYR
41	BG	3	LEU
41	BG	48	GLU
41	BG	53	LEU
41	BG	81	LYS
41	BG	82	LEU
41	BG	84	LYS
41	BG	86	MET
41	BG	87	PRO
41	BG	90	LEU
41	BG	110	ALA
41	BG	145	THR

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Mol	Chain	Res	Type
41	BG	151	ALA
41	BG	171	ALA
41	BG	172	LEU
42	BH	13	LYS
42	BH	44	VAL
42	BH	71	LEU
42	BH	85	LYS
42	BH	89	ILE
42	BH	92	ILE
42	BH	137	ASP
42	BH	138	LYS
42	BH	151	ILE
42	BH	155	SER
42	BH	156	ALA
42	BH	159	GLU
42	BH	160	LYS
42	BH	165	ALA
43	BI	15	VAL
43	BI	81	VAL
43	BI	86	THR
43	BI	88	ILE
43	BI	91	SER
43	BI	133	HIS
43	BI	145	VAL
44	BN	8	GLN
44	BN	44	PRO
44	BN	47	ALA
44	BN	52	VAL
44	BN	58	ASP
44	BN	74	ARG
44	BN	78	TYR
44	BN	127	ASP
44	BN	130	HIS
44	BN	133	GLN
45	BO	14	THR
45	BO	35	VAL
45	BO	48	PRO
45	BO	102	VAL
45	BO	112	MET
46	BP	15	ARG
46	BP	17	LYS
46	BP	18	ARG

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Mol	Chain	Res	Type
46	BP	31	ALA
46	BP	35	HIS
46	BP	46	LYS
46	BP	49	ARG
46	BP	64	LYS
46	BP	65	ARG
46	BP	67	MET
46	BP	111	ARG
46	BP	146	VAL
46	BP	147	LEU
47	BQ	7	MET
47	BQ	11	LYS
47	BQ	12	GLN
47	BQ	25	ASP
47	BQ	66	ILE
47	BQ	71	ASP
47	BQ	88	GLY
47	BQ	103	MET
47	BQ	134	ARG
47	BQ	136	ALA
47	BQ	138	ASP
48	BR	7	GLY
48	BR	8	ARG
48	BR	12	ARG
48	BR	14	SER
48	BR	16	HIS
48	BR	29	LEU
48	BR	45	ARG
48	BR	71	GLN
48	BR	86	ARG
48	BR	117	VAL
49	BS	17	ARG
49	BS	24	LEU
49	BS	33	LYS
49	BS	42	ASP
49	BS	56	LEU
49	BS	59	LYS
49	BS	74	ALA
49	BS	83	LYS
49	BS	84	GLN
49	BS	92	TYR
49	BS	93	LYS

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Mol	Chain	Res	Type
49	BS	96	GLY
49	BS	100	ALA
49	BS	102	ALA
49	BS	105	ALA
50	BT	18	ASP
50	BT	20	PRO
50	BT	24	PRO
50	BT	28	VAL
50	BT	30	VAL
50	BT	32	TYR
50	BT	33	LYS
50	BT	41	ARG
50	BT	57	PHE
50	BT	80	SER
50	BT	83	ILE
50	BT	97	ALA
50	BT	107	ASP
51	BU	5	LYS
51	BU	7	GLY
51	BU	8	VAL
51	BU	9	VAL
51	BU	45	TYR
51	BU	53	ARG
51	BU	93	LYS
51	BU	97	ASP
52	BV	19	LYS
52	BV	23	GLU
52	BV	24	LYS
52	BV	28	GLU
52	BV	50	PRO
52	BV	51	VAL
52	BV	57	VAL
52	BV	59	ALA
52	BV	73	SER
52	BV	98	GLU
53	BW	11	ARG
53	BW	63	ASP
53	BW	75	TYR
53	BW	93	ALA
53	BW	111	HIS
54	BX	6	ASP
54	BX	26	TYR

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Mol	Chain	Res	Type
54	BX	36	LYS
54	BX	53	LYS
54	BX	59	VAL
54	BX	60	ARG
54	BX	62	LYS
54	BX	65	ARG
54	BX	73	ARG
54	BX	84	ALA
55	BY	3	VAL
55	BY	7	VAL
55	BY	16	ALA
55	BY	38	ILE
55	BY	44	ILE
55	BY	55	TYR
55	BY	56	PRO
55	BY	66	PRO
55	BY	74	PRO
55	BY	75	ILE
55	BY	76	CYS
55	BY	78	ALA
55	BY	90	LEU
55	BY	96	ILE
55	BY	101	LYS
56	BZ	8	TYR
56	BZ	30	ASN
56	BZ	48	PHE
56	BZ	77	ASP
56	BZ	78	LYS
56	BZ	81	ARG
56	BZ	93	ASP
56	BZ	100	VAL
56	BZ	111	VAL
56	BZ	113	ALA
56	BZ	128	VAL
56	BZ	135	GLU
56	BZ	149	SER
56	BZ	153	SER
56	BZ	163	LEU
56	BZ	166	SER
56	BZ	168	GLU
2	CB	20	GLU
2	CB	22	LYS

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Mol	Chain	Res	Type
2	CB	23	ARG
2	CB	29	ALA
2	CB	30	ARG
2	CB	37	ASN
2	CB	75	LYS
2	CB	77	ALA
2	CB	88	ALA
2	CB	154	LEU
2	CB	165	VAL
2	CB	167	PRO
2	CB	168	THR
2	CB	191	ASP
2	CB	194	PRO
2	CB	202	PRO
2	CB	204	ASN
2	CB	209	ARG
2	CB	229	VAL
2	CB	238	LEU
2	CB	239	VAL
3	CC	4	LYS
3	CC	12	LEU
3	CC	53	ALA
3	CC	61	ALA
3	CC	157	ILE
3	CC	172	ARG
3	CC	174	PRO
3	CC	191	THR
3	CC	207	VAL
4	CD	4	TYR
4	CD	5	ILE
4	CD	6	GLY
4	CD	9	CYS
4	CD	14	ARG
4	CD	24	GLU
4	CD	28	SER
4	CD	119	GLN
4	CD	128	VAL
4	CD	149	ALA
4	CD	150	GLU
4	CD	163	GLU
4	CD	164	ALA
4	CD	177	ASP

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Mol	Chain	Res	Type
4	CD	191	ARG
5	CE	37	ARG
5	CE	129	ILE
6	CF	13	ASN
6	CF	69	GLU
6	CF	70	ASP
6	CF	72	VAL
7	CG	31	MET
7	CG	39	ALA
7	CG	153	HIS
8	CH	18	ARG
8	CH	22	GLU
8	CH	29	SER
8	CH	97	VAL
8	CH	129	VAL
9	CI	12	GLU
9	CI	29	ASN
9	CI	117	HIS
9	CI	120	ARG
10	CJ	33	GLN
10	CJ	57	LYS
10	CJ	91	PRO
11	CK	27	ASN
11	CK	103	LEU
11	CK	105	VAL
11	CK	106	LYS
12	CL	10	LEU
12	CL	23	LYS
12	CL	28	LYS
12	CL	47	LYS
12	CL	48	PRO
12	CL	62	SER
13	CM	4	ILE
13	CM	11	ARG
13	CM	60	VAL
13	CM	68	GLY
13	CM	90	LEU
13	CM	117	VAL
14	CN	22	THR
14	CN	23	ARG
15	CO	16	ALA
17	CQ	49	GLU

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Mol	Chain	Res	Type
17	CQ	74	LEU
17	CQ	93	GLN
18	CR	20	ALA
18	CR	37	VAL
18	CR	45	SER
19	CS	10	PHE
19	CS	67	VAL
19	CS	80	TYR
19	CS	81	ARG
20	CT	28	ALA
20	CT	49	ALA
20	CT	63	ILE
20	CT	64	ASP
25	CY	19	GLU
25	CY	30	THR
25	CY	52	LEU
25	CY	53	ASN
25	CY	87	ASP
25	CY	120	GLN
25	CY	132	ILE
25	CY	150	SER
25	CY	166	ASP
26	D0	17	GLN
26	D0	55	ARG
26	D0	84	LEU
27	D1	12	PRO
27	D1	15	ALA
27	D1	26	ARG
27	D1	33	LYS
27	D1	53	VAL
27	D1	57	GLU
27	D1	63	ALA
28	D2	14	ARG
28	D2	15	LYS
28	D2	28	LYS
28	D2	50	ILE
28	D2	51	ARG
29	D3	3	ARG
29	D3	13	ILE
29	D3	31	LEU
29	D3	51	ALA
29	D3	52	HIS

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Mol	Chain	Res	Type
29	D3	56	VAL
29	D3	57	GLU
30	D4	6	HIS
30	D4	7	PRO
30	D4	10	VAL
30	D4	13	ARG
30	D4	14	ILE
30	D4	23	GLU
30	D4	29	PRO
30	D4	35	VAL
30	D4	47	GLN
31	D5	21	SER
31	D5	22	HIS
31	D5	32	PRO
31	D5	33	CYS
31	D5	34	PRO
31	D5	35	GLU
31	D5	57	VAL
32	D6	16	CYS
32	D6	31	PRO
32	D6	44	ARG
32	D6	45	LYS
33	D7	18	PHE
33	D7	46	VAL
34	D8	4	MET
34	D8	25	MET
34	D8	32	LEU
34	D8	37	SER
34	D8	43	GLN
37	DC	22	ILE
37	DC	47	LEU
37	DC	49	ILE
37	DC	54	SER
37	DC	68	LEU
37	DC	104	LEU
37	DC	126	LYS
37	DC	133	PRO
37	DC	140	PRO
37	DC	167	LYS
37	DC	173	ALA
37	DC	174	PRO
37	DC	184	LYS

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Mol	Chain	Res	Type
37	DC	204	ALA
37	DC	222	VAL
38	DD	3	VAL
38	DD	8	PRO
38	DD	13	ARG
38	DD	22	SER
38	DD	25	THR
38	DD	28	GLU
38	DD	32	SER
38	DD	45	ASN
38	DD	53	PHE
38	DD	58	HIS
38	DD	135	PHE
38	DD	159	ALA
38	DD	204	ILE
38	DD	216	GLY
38	DD	223	GLY
38	DD	224	ALA
38	DD	228	PRO
38	DD	241	PRO
38	DD	262	ARG
38	DD	263	ARG
38	DD	266	SER
38	DD	272	ALA
39	DE	2	LYS
39	DE	4	ILE
39	DE	45	THR
39	DE	53	PRO
39	DE	57	LYS
39	DE	61	ARG
39	DE	66	HIS
39	DE	77	ILE
39	DE	82	ARG
39	DE	88	GLY
39	DE	90	THR
39	DE	93	VAL
39	DE	122	PHE
39	DE	129	HIS
39	DE	182	LEU
39	DE	186	GLY
40	DF	2	LYS
40	DF	6	VAL

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Mol	Chain	Res	Type
40	DF	10	PRO
40	DF	21	ALA
40	DF	40	GLN
40	DF	57	VAL
40	DF	60	SER
40	DF	66	PRO
40	DF	72	ARG
40	DF	73	ALA
40	DF	85	GLY
40	DF	89	VAL
40	DF	184	TYR
41	DG	3	LEU
41	DG	6	ALA
41	DG	12	TYR
41	DG	48	GLU
41	DG	50	ALA
41	DG	53	LEU
41	DG	77	ILE
41	DG	82	LEU
41	DG	84	LYS
41	DG	86	MET
41	DG	104	GLU
41	DG	111	LEU
41	DG	122	PRO
41	DG	123	ASN
41	DG	130	ASN
41	DG	142	PRO
41	DG	143	GLU
41	DG	150	ASP
41	DG	157	ILE
41	DG	163	ALA
41	DG	172	LEU
42	DH	13	LYS
42	DH	44	VAL
42	DH	71	LEU
42	DH	85	LYS
42	DH	89	ILE
42	DH	92	ILE
42	DH	137	ASP
42	DH	138	LYS
42	DH	151	ILE
42	DH	155	SER

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Mol	Chain	Res	Type
42	DH	156	ALA
42	DH	159	GLU
42	DH	160	LYS
42	DH	165	ALA
43	DI	15	VAL
43	DI	81	VAL
43	DI	86	THR
43	DI	88	ILE
43	DI	91	SER
43	DI	133	HIS
43	DI	145	VAL
44	DN	8	GLN
44	DN	44	PRO
44	DN	47	ALA
44	DN	52	VAL
44	DN	58	ASP
44	DN	74	ARG
44	DN	78	TYR
44	DN	127	ASP
44	DN	130	HIS
44	DN	133	GLN
45	DO	14	THR
45	DO	35	VAL
45	DO	48	PRO
45	DO	102	VAL
45	DO	112	MET
46	DP	15	ARG
46	DP	17	LYS
46	DP	18	ARG
46	DP	31	ALA
46	DP	35	HIS
46	DP	46	LYS
46	DP	49	ARG
46	DP	64	LYS
46	DP	65	ARG
46	DP	67	MET
46	DP	111	ARG
46	DP	146	VAL
46	DP	147	LEU
47	DQ	7	MET
47	DQ	11	LYS
47	DQ	12	GLN

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Mol	Chain	Res	Type
47	DQ	25	ASP
47	DQ	66	ILE
47	DQ	71	ASP
47	DQ	88	GLY
47	DQ	103	MET
47	DQ	134	ARG
47	DQ	136	ALA
47	DQ	138	ASP
48	DR	7	GLY
48	DR	8	ARG
48	DR	12	ARG
48	DR	14	SER
48	DR	16	HIS
48	DR	29	LEU
48	DR	45	ARG
48	DR	71	GLN
48	DR	86	ARG
48	DR	117	VAL
49	DS	17	ARG
49	DS	24	LEU
49	DS	33	LYS
49	DS	42	ASP
49	DS	56	LEU
49	DS	59	LYS
49	DS	74	ALA
49	DS	83	LYS
49	DS	84	GLN
49	DS	92	TYR
49	DS	93	LYS
49	DS	96	GLY
49	DS	100	ALA
49	DS	102	ALA
49	DS	105	ALA
50	DT	18	ASP
50	DT	20	PRO
50	DT	24	PRO
50	DT	28	VAL
50	DT	30	VAL
50	DT	32	TYR
50	DT	33	LYS
50	DT	41	ARG
50	DT	57	PHE

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Mol	Chain	Res	Type
50	DT	80	SER
50	DT	83	ILE
50	DT	97	ALA
50	DT	107	ASP
51	DU	5	LYS
51	DU	7	GLY
51	DU	8	VAL
51	DU	9	VAL
51	DU	22	LYS
51	DU	45	TYR
51	DU	53	ARG
51	DU	93	LYS
51	DU	97	ASP
52	DV	19	LYS
52	DV	23	GLU
52	DV	24	LYS
52	DV	28	GLU
52	DV	50	PRO
52	DV	51	VAL
52	DV	57	VAL
52	DV	59	ALA
52	DV	73	SER
52	DV	98	GLU
53	DW	11	ARG
53	DW	63	ASP
53	DW	75	TYR
53	DW	93	ALA
53	DW	111	HIS
54	DX	6	ASP
54	DX	26	TYR
54	DX	37	THR
54	DX	53	LYS
54	DX	59	VAL
54	DX	60	ARG
54	DX	62	LYS
54	DX	65	ARG
54	DX	73	ARG
54	DX	84	ALA
55	DY	3	VAL
55	DY	7	VAL
55	DY	16	ALA
55	DY	38	ILE

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Mol	Chain	Res	Type
55	DY	44	ILE
55	DY	55	TYR
55	DY	56	PRO
55	DY	64	GLU
55	DY	66	PRO
55	DY	74	PRO
55	DY	75	ILE
55	DY	76	CYS
55	DY	78	ALA
55	DY	90	LEU
55	DY	96	ILE
55	DY	101	LYS
56	DZ	30	ASN
56	DZ	33	LEU
56	DZ	78	LYS
56	DZ	101	PRO
56	DZ	111	VAL
56	DZ	135	GLU
56	DZ	142	SER
56	DZ	148	ASP
56	DZ	149	SER
56	DZ	152	ALA
56	DZ	166	SER
56	DZ	169	GLU
2	AB	18	GLY
2	AB	23	ARG
2	AB	32	ILE
2	AB	45	GLN
2	AB	52	GLU
2	AB	64	ARG
2	AB	78	GLN
2	AB	84	GLU
2	AB	100	GLY
2	AB	105	PHE
2	AB	160	ASP
2	AB	204	ASN
2	AB	235	SER
3	AC	5	ILE
3	AC	18	TRP
3	AC	38	ARG
3	AC	47	LEU
3	AC	52	LEU

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Mol	Chain	Res	Type
3	AC	54	ARG
3	AC	74	GLY
3	AC	127	ARG
3	AC	164	ARG
3	AC	181	ASN
3	AC	189	ALA
4	AD	3	ARG
4	AD	15	GLU
4	AD	30	LYS
4	AD	44	GLY
4	AD	47	ARG
4	AD	57	ARG
4	AD	60	GLU
4	AD	89	THR
4	AD	132	ARG
4	AD	159	ARG
5	AE	27	ARG
5	AE	38	GLN
5	AE	85	GLY
5	AE	86	ALA
5	AE	113	ALA
5	AE	130	ASN
5	AE	146	ALA
6	AF	62	TRP
6	AF	80	ARG
6	AF	84	ASN
7	AG	7	ALA
7	AG	105	VAL
8	AH	64	LYS
9	AI	12	GLU
9	AI	40	LEU
9	AI	100	GLY
10	AJ	23	ILE
10	AJ	36	GLY
10	AJ	59	SER
10	AJ	84	GLN
10	AJ	86	MET
11	AK	16	SER
11	AK	25	TYR
11	AK	103	LEU
11	AK	106	LYS
12	AL	56	ALA

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Mol	Chain	Res	Type
12	AL	83	VAL
12	AL	91	LYS
12	AL	96	VAL
12	AL	102	ARG
13	AM	67	GLU
13	AM	100	GLY
13	AM	107	ALA
13	AM	124	PRO
14	AN	16	PHE
14	AN	52	GLN
15	AO	26	GLU
15	AO	65	ARG
15	AO	73	GLU
16	AP	28	ARG
16	AP	53	VAL
16	AP	63	GLY
17	AQ	33	GLY
17	AQ	80	GLY
18	AR	31	LEU
18	AR	59	SER
18	AR	85	LEU
18	AR	86	VAL
19	AS	26	GLY
19	AS	53	ASN
19	AS	70	LYS
20	AT	25	ARG
20	AT	42	GLN
20	AT	78	ALA
20	AT	99	LEU
20	AT	102	GLY
21	AU	9	ARG
25	AY	30	THR
25	AY	71	TRP
26	B0	4	LYS
26	B0	11	ARG
26	B0	20	ARG
27	B1	10	LYS
27	B1	28	GLY
27	B1	47	GLN
27	B1	64	ALA
27	B1	66	HIS
27	B1	71	TYR

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Mol	Chain	Res	Type
27	B1	82	LEU
28	B2	15	LYS
28	B2	36	ARG
28	B2	41	ILE
28	B2	57	ILE
28	B2	60	LEU
29	B3	27	GLY
30	B4	3	GLU
30	B4	16	CYS
30	B4	31	ILE
31	B5	12	SER
31	B5	49	CYS
31	B5	51	TYR
32	B6	49	HIS
32	B6	52	VAL
33	B7	36	GLN
34	B8	7	HIS
34	B8	35	GLN
37	BC	35	ALA
37	BC	36	LYS
37	BC	77	ILE
37	BC	78	ALA
37	BC	106	ALA
37	BC	134	ARG
37	BC	135	GLY
37	BC	141	LYS
37	BC	147	PHE
37	BC	171	ILE
37	BC	215	THR
37	BC	217	THR
38	BD	26	LYS
38	BD	53	PHE
38	BD	57	GLY
38	BD	70	TRP
38	BD	125	ILE
38	BD	134	ARG
38	BD	156	ALA
38	BD	202	LYS
38	BD	234	GLY
38	BD	239	ARG
38	BD	265	PRO
38	BD	268	ARG

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Mol	Chain	Res	Type
39	BE	30	PRO
39	BE	54	GLN
39	BE	56	PRO
39	BE	60	ASN
39	BE	73	GLU
39	BE	84	PHE
39	BE	92	THR
39	BE	121	ASN
39	BE	144	ARG
39	BE	164	ARG
39	BE	169	ASN
39	BE	174	ASP
39	BE	185	LYS
39	BE	189	PRO
39	BE	191	PRO
40	BF	22	ALA
40	BF	59	TYR
40	BF	64	ILE
40	BF	68	LYS
40	BF	69	HIS
40	BF	70	THR
40	BF	86	GLY
40	BF	103	LYS
40	BF	111	ALA
40	BF	122	LYS
40	BF	127	GLU
40	BF	134	GLY
40	BF	165	ARG
40	BF	172	TRP
40	BF	178	PRO
41	BG	5	VAL
41	BG	14	GLU
41	BG	45	GLU
41	BG	61	ALA
41	BG	68	PRO
41	BG	96	ARG
41	BG	127	GLY
41	BG	129	GLY
41	BG	143	GLU
42	BH	14	GLY
42	BH	16	SER
42	BH	90	LYS

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Mol	Chain	Res	Type
42	BH	100	GLY
42	BH	141	VAL
42	BH	146	ALA
42	BH	147	ASN
42	BH	148	ILE
43	BI	74	ASN
43	BI	82	ARG
43	BI	111	PRO
43	BI	112	LYS
44	BN	4	TYR
44	BN	9	VAL
44	BN	42	TRP
44	BN	45	ASN
44	BN	60	ILE
44	BN	61	ARG
44	BN	64	GLY
44	BN	83	LYS
44	BN	98	VAL
44	BN	108	PRO
44	BN	109	LYS
45	BO	27	GLY
45	BO	44	LYS
45	BO	80	ASP
45	BO	81	ASP
46	BP	34	GLY
46	BP	38	GLN
46	BP	39	LYS
46	BP	42	SER
46	BP	56	SER
46	BP	66	GLY
46	BP	102	ARG
46	BP	107	LYS
46	BP	108	LYS
47	BQ	62	GLY
48	BR	9	LYS
48	BR	10	LEU
48	BR	42	LYS
48	BR	49	ASP
48	BR	82	GLU
48	BR	105	ARG
48	BR	107	ASP
49	BS	32	LEU

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Mol	Chain	Res	Type
49	BS	62	LYS
49	BS	85	VAL
49	BS	89	ARG
49	BS	94	TYR
49	BS	98	VAL
50	BT	52	ILE
50	BT	81	PRO
50	BT	84	GLN
50	BT	92	GLY
50	BT	101	PHE
50	BT	115	ARG
50	BT	136	GLN
51	BU	22	LYS
51	BU	24	TYR
51	BU	32	PHE
51	BU	60	LEU
51	BU	62	ILE
51	BU	65	ILE
51	BU	79	PHE
52	BV	17	GLY
52	BV	36	PRO
52	BV	41	GLY
52	BV	44	LYS
52	BV	46	VAL
52	BV	47	VAL
52	BV	64	HIS
52	BV	68	LYS
52	BV	69	LYS
52	BV	79	VAL
52	BV	86	GLY
52	BV	91	TYR
53	BW	6	ILE
53	BW	18	ARG
53	BW	25	ARG
53	BW	35	ILE
53	BW	60	ASN
54	BX	7	VAL
54	BX	24	GLY
54	BX	37	THR
54	BX	49	VAL
54	BX	68	ARG
54	BX	69	TYR

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Mol	Chain	Res	Type
54	BX	76	ARG
54	BX	81	VAL
54	BX	88	LYS
55	BY	29	GLU
55	BY	64	GLU
55	BY	67	LEU
56	BZ	41	LEU
56	BZ	52	SER
56	BZ	62	PRO
56	BZ	63	ASP
56	BZ	66	SER
56	BZ	98	MET
56	BZ	101	PRO
56	BZ	109	ALA
56	BZ	115	GLY
56	BZ	138	GLU
56	BZ	140	ASP
56	BZ	147	GLY
56	BZ	148	ASP
56	BZ	165	VAL
56	BZ	177	PRO
2	CB	18	GLY
2	CB	32	ILE
2	CB	45	GLN
2	CB	52	GLU
2	CB	64	ARG
2	CB	78	GLN
2	CB	84	GLU
2	CB	100	GLY
2	CB	160	ASP
2	CB	235	SER
3	CC	5	ILE
3	CC	18	TRP
3	CC	47	LEU
3	CC	52	LEU
3	CC	54	ARG
3	CC	74	GLY
3	CC	127	ARG
3	CC	160	ALA
3	CC	164	ARG
3	CC	181	ASN
4	CD	3	ARG

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Mol	Chain	Res	Type
4	CD	10	ARG
4	CD	15	GLU
4	CD	30	LYS
4	CD	44	GLY
4	CD	47	ARG
4	CD	57	ARG
4	CD	59	ARG
4	CD	60	GLU
4	CD	89	THR
4	CD	97	LEU
4	CD	132	ARG
4	CD	159	ARG
4	CD	193	ASP
5	CE	27	ARG
5	CE	49	PRO
5	CE	85	GLY
5	CE	86	ALA
5	CE	130	ASN
5	CE	146	ALA
6	CF	62	TRP
6	CF	80	ARG
6	CF	84	ASN
7	CG	7	ALA
7	CG	105	VAL
7	CG	120	ILE
7	CG	149	ARG
8	CH	17	THR
8	CH	64	LYS
9	CI	40	LEU
9	CI	100	GLY
10	CJ	23	ILE
10	CJ	36	GLY
10	CJ	59	SER
10	CJ	84	GLN
10	CJ	86	MET
11	CK	16	SER
11	CK	25	TYR
11	CK	26	ASN
11	CK	63	LEU
12	CL	56	ALA
12	CL	83	VAL
12	CL	91	LYS

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Mol	Chain	Res	Type
12	CL	96	VAL
12	CL	102	ARG
12	CL	123	LYS
13	CM	83	ASP
13	CM	100	GLY
13	CM	107	ALA
13	CM	124	PRO
14	CN	16	PHE
14	CN	52	GLN
15	CO	26	GLU
15	CO	65	ARG
15	CO	73	GLU
16	CP	28	ARG
16	CP	53	VAL
16	CP	63	GLY
17	CQ	4	LYS
17	CQ	33	GLY
17	CQ	80	GLY
17	CQ	81	ARG
17	CQ	96	GLU
18	CR	31	LEU
18	CR	41	LYS
18	CR	59	SER
18	CR	85	LEU
18	CR	86	VAL
19	CS	26	GLY
19	CS	53	ASN
19	CS	70	LYS
20	CT	25	ARG
20	CT	42	GLN
20	CT	78	ALA
20	CT	99	LEU
20	CT	102	GLY
25	CY	23	HIS
25	CY	24	ASN
25	CY	29	ARG
25	CY	39	LEU
25	CY	63	PRO
25	CY	64	ARG
25	CY	83	ILE
25	CY	107	THR
25	CY	109	GLU

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Mol	Chain	Res	Type
25	CY	126	ARG
25	CY	127	VAL
25	CY	161	ILE
25	CY	165	THR
25	CY	174	GLN
26	D0	4	LYS
26	D0	11	ARG
26	D0	20	ARG
27	D1	13	ILE
27	D1	34	THR
27	D1	38	SER
27	D1	49	VAL
27	D1	56	GLN
27	D1	68	PRO
27	D1	69	LYS
27	D1	77	ALA
27	D1	79	GLY
27	D1	82	LEU
27	D1	87	PRO
27	D1	89	GLU
28	D2	21	LEU
28	D2	40	SER
28	D2	41	ILE
28	D2	53	LEU
28	D2	60	LEU
29	D3	27	GLY
30	D4	3	GLU
30	D4	16	CYS
30	D4	31	ILE
31	D5	12	SER
31	D5	49	CYS
31	D5	51	TYR
32	D6	49	HIS
32	D6	52	VAL
33	D7	36	GLN
34	D8	7	HIS
34	D8	35	GLN
37	DC	35	ALA
37	DC	36	LYS
37	DC	52	ARG
37	DC	77	ILE
37	DC	78	ALA

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Mol	Chain	Res	Type
37	DC	106	ALA
37	DC	125	SER
37	DC	134	ARG
37	DC	135	GLY
37	DC	141	LYS
37	DC	147	PHE
37	DC	171	ILE
37	DC	215	THR
37	DC	217	THR
38	DD	26	LYS
38	DD	57	GLY
38	DD	70	TRP
38	DD	125	ILE
38	DD	134	ARG
38	DD	156	ALA
38	DD	200	ASP
38	DD	202	LYS
38	DD	206	LEU
38	DD	234	GLY
38	DD	239	ARG
38	DD	265	PRO
38	DD	268	ARG
39	DE	30	PRO
39	DE	54	GLN
39	DE	56	PRO
39	DE	73	GLU
39	DE	84	PHE
39	DE	89	ASP
39	DE	92	THR
39	DE	121	ASN
39	DE	144	ARG
39	DE	164	ARG
39	DE	169	ASN
39	DE	174	ASP
39	DE	185	LYS
39	DE	189	PRO
39	DE	191	PRO
40	DF	22	ALA
40	DF	59	TYR
40	DF	68	LYS
40	DF	69	HIS
40	DF	70	THR

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Mol	Chain	Res	Type
40	DF	86	GLY
40	DF	103	LYS
40	DF	111	ALA
40	DF	122	LYS
40	DF	127	GLU
40	DF	134	GLY
40	DF	165	ARG
40	DF	172	TRP
40	DF	178	PRO
41	DG	5	VAL
41	DG	46	ALA
41	DG	81	LYS
41	DG	105	LYS
41	DG	107	LEU
41	DG	129	GLY
41	DG	139	LEU
41	DG	151	ALA
41	DG	164	GLU
42	DH	14	GLY
42	DH	16	SER
42	DH	90	LYS
42	DH	100	GLY
42	DH	141	VAL
42	DH	146	ALA
42	DH	147	ASN
43	DI	74	ASN
43	DI	82	ARG
43	DI	111	PRO
43	DI	112	LYS
44	DN	4	TYR
44	DN	9	VAL
44	DN	14	VAL
44	DN	42	TRP
44	DN	45	ASN
44	DN	60	ILE
44	DN	61	ARG
44	DN	64	GLY
44	DN	83	LYS
44	DN	98	VAL
44	DN	108	PRO
44	DN	109	LYS
45	DO	44	LYS

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Mol	Chain	Res	Type
45	DO	80	ASP
45	DO	81	ASP
46	DP	34	GLY
46	DP	37	GLY
46	DP	38	GLN
46	DP	39	LYS
46	DP	42	SER
46	DP	56	SER
46	DP	66	GLY
46	DP	102	ARG
46	DP	107	LYS
46	DP	108	LYS
46	DP	122	PRO
47	DQ	62	GLY
47	DQ	90	VAL
48	DR	9	LYS
48	DR	10	LEU
48	DR	15	SER
48	DR	42	LYS
48	DR	49	ASP
48	DR	68	ARG
48	DR	82	GLU
49	DS	32	LEU
49	DS	62	LYS
49	DS	89	ARG
49	DS	94	TYR
49	DS	98	VAL
50	DT	52	ILE
50	DT	81	PRO
50	DT	84	GLN
50	DT	92	GLY
50	DT	101	PHE
50	DT	115	ARG
50	DT	136	GLN
51	DU	24	TYR
51	DU	32	PHE
51	DU	60	LEU
51	DU	62	ILE
51	DU	65	ILE
51	DU	79	PHE
52	DV	17	GLY
52	DV	36	PRO

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Mol	Chain	Res	Type
52	DV	41	GLY
52	DV	44	LYS
52	DV	46	VAL
52	DV	47	VAL
52	DV	64	HIS
52	DV	68	LYS
52	DV	69	LYS
52	DV	72	VAL
52	DV	79	VAL
52	DV	86	GLY
52	DV	91	TYR
53	DW	6	ILE
53	DW	25	ARG
53	DW	35	ILE
53	DW	60	ASN
54	DX	24	GLY
54	DX	36	LYS
54	DX	49	VAL
54	DX	68	ARG
54	DX	69	TYR
54	DX	76	ARG
54	DX	81	VAL
54	DX	82	GLN
54	DX	88	LYS
55	DY	67	LEU
56	DZ	45	ASP
56	DZ	65	GLN
56	DZ	108	PRO
56	DZ	115	GLY
56	DZ	121	HIS
56	DZ	147	GLY
56	DZ	163	LEU
56	DZ	171	ILE
2	AB	83	MET
2	AB	106	LYS
2	AB	133	LYS
2	AB	135	GLN
2	AB	143	GLU
2	AB	176	GLU
2	AB	178	ARG
2	AB	195	ASP
2	AB	213	LEU

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Mol	Chain	Res	Type
3	AC	60	ALA
3	AC	107	GLN
3	AC	117	ALA
3	AC	145	GLY
3	AC	160	ALA
4	AD	10	ARG
4	AD	29	PRO
4	AD	43	HIS
4	AD	59	ARG
4	AD	64	LEU
4	AD	97	LEU
4	AD	101	LEU
4	AD	147	ALA
4	AD	153	ARG
4	AD	193	ASP
5	AE	49	PRO
5	AE	73	ASN
5	AE	96	PRO
5	AE	132	ALA
6	AF	17	SER
6	AF	29	ALA
6	AF	71	ARG
7	AG	14	PRO
7	AG	40	ALA
7	AG	65	ALA
7	AG	100	ALA
7	AG	120	ILE
7	AG	149	ARG
8	AH	3	THR
8	AH	6	ILE
8	AH	37	ARG
8	AH	41	ARG
8	AH	68	ARG
8	AH	121	ASP
9	AI	94	ALA
11	AK	26	ASN
11	AK	63	LEU
11	AK	100	ALA
11	AK	124	LYS
11	AK	126	ARG
12	AL	46	LYS
12	AL	79	GLU

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Mol	Chain	Res	Type
12	AL	123	LYS
13	AM	61	GLU
13	AM	102	ARG
14	AN	5	ALA
14	AN	35	ARG
15	AO	13	GLN
15	AO	37	ASN
16	AP	13	HIS
16	AP	73	LEU
16	AP	81	ARG
16	AP	83	GLU
17	AQ	4	LYS
17	AQ	78	GLU
17	AQ	87	LYS
18	AR	41	LYS
18	AR	43	PHE
18	AR	82	THR
19	AS	44	MET
20	AT	12	ALA
20	AT	95	ALA
20	AT	96	GLY
20	AT	97	ALA
21	AU	12	LYS
25	AY	39	LEU
25	AY	53	ASN
25	AY	64	ARG
25	AY	87	ASP
25	AY	88	LEU
25	AY	92	PRO
25	AY	143	LEU
25	AY	184	LEU
26	B0	2	ALA
27	B1	56	GLN
29	B3	29	ARG
30	B4	44	THR
31	B5	58	LEU
33	B7	19	ARG
33	B7	32	LYS
34	B8	13	ARG
34	B8	36	LYS
34	B8	51	ALA
34	B8	64	TYR

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Mol	Chain	Res	Type
37	BC	24	GLU
37	BC	51	PRO
37	BC	52	ARG
37	BC	125	SER
37	BC	146	GLY
37	BC	153	ILE
37	BC	172	HIS
37	BC	183	GLU
37	BC	201	PRO
37	BC	211	SER
37	BC	214	VAL
38	BD	89	SER
38	BD	115	GLN
38	BD	200	ASP
38	BD	206	LEU
38	BD	214	TRP
39	BE	43	GLY
39	BE	44	TYR
39	BE	89	ASP
39	BE	162	ALA
39	BE	180	ASN
40	BF	14	PRO
40	BF	33	LEU
40	BF	61	GLY
40	BF	115	ALA
40	BF	132	VAL
40	BF	133	ASN
40	BF	179	GLU
41	BG	25	TYR
41	BG	43	LEU
41	BG	117	PHE
41	BG	157	ILE
41	BG	175	LEU
42	BH	39	PRO
42	BH	152	ARG
42	BH	153	LYS
42	BH	154	PRO
43	BI	6	LEU
43	BI	40	THR
43	BI	85	GLU
43	BI	104	GLN
43	BI	120	ILE

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Mol	Chain	Res	Type
44	BN	14	VAL
44	BN	18	ALA
44	BN	46	VAL
44	BN	57	ALA
44	BN	80	GLY
44	BN	129	PRO
45	BO	36	GLY
45	BO	45	GLU
45	BO	68	GLU
45	BO	90	GLN
45	BO	114	ILE
46	BP	12	ALA
46	BP	33	ARG
46	BP	110	TYR
46	BP	120	ALA
46	BP	122	PRO
46	BP	129	ALA
46	BP	141	ALA
47	BQ	24	GLY
47	BQ	59	ARG
47	BQ	60	ARG
47	BQ	78	PRO
47	BQ	90	VAL
47	BQ	105	GLU
47	BQ	115	MET
47	BQ	140	ALA
48	BR	5	LYS
48	BR	15	SER
48	BR	17	ARG
48	BR	50	HIS
48	BR	60	LEU
48	BR	68	ARG
48	BR	88	ARG
49	BS	43	GLU
49	BS	53	SER
49	BS	54	LEU
49	BS	61	ASN
49	BS	73	LEU
49	BS	78	LEU
49	BS	97	ARG
50	BT	23	ARG
50	BT	25	GLY

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Mol	Chain	Res	Type
50	BT	131	ALA
51	BU	83	LEU
51	BU	88	ILE
51	BU	91	ASP
51	BU	92	ARG
52	BV	40	LEU
52	BV	42	GLY
52	BV	43	GLU
52	BV	48	GLY
52	BV	53	GLU
52	BV	72	VAL
52	BV	82	ARG
53	BW	14	PRO
53	BW	66	GLU
54	BX	82	GLN
54	BX	89	ILE
55	BY	23	ARG
55	BY	31	LEU
55	BY	39	VAL
55	BY	40	GLU
55	BY	99	CYS
56	BZ	83	PRO
56	BZ	126	VAL
56	BZ	130	PRO
56	BZ	141	VAL
2	CB	83	MET
2	CB	105	PHE
2	CB	106	LYS
2	CB	133	LYS
2	CB	135	GLN
2	CB	143	GLU
2	CB	176	GLU
2	CB	178	ARG
2	CB	195	ASP
2	CB	213	LEU
3	CC	38	ARG
3	CC	60	ALA
3	CC	107	GLN
3	CC	117	ALA
3	CC	145	GLY
3	CC	179	ARG
3	CC	189	ALA

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Mol	Chain	Res	Type
4	CD	29	PRO
4	CD	64	LEU
4	CD	101	LEU
4	CD	129	ASN
4	CD	147	ALA
5	CE	38	GLN
5	CE	73	ASN
5	CE	96	PRO
5	CE	113	ALA
5	CE	132	ALA
6	CF	17	SER
6	CF	29	ALA
6	CF	71	ARG
7	CG	14	PRO
7	CG	40	ALA
7	CG	65	ALA
7	CG	100	ALA
7	CG	119	ARG
7	CG	131	LYS
8	CH	3	THR
8	CH	6	ILE
8	CH	37	ARG
8	CH	41	ARG
8	CH	121	ASP
9	CI	94	ALA
11	CK	18	ARG
11	CK	66	LEU
11	CK	100	ALA
11	CK	124	LYS
11	CK	126	ARG
12	CL	79	GLU
13	CM	61	GLU
13	CM	67	GLU
13	CM	102	ARG
14	CN	5	ALA
14	CN	35	ARG
15	CO	13	GLN
16	CP	13	HIS
16	CP	73	LEU
16	CP	81	ARG
17	CQ	17	LYS
17	CQ	78	GLU

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Mol	Chain	Res	Type
17	CQ	87	LYS
18	CR	43	PHE
18	CR	82	THR
19	CS	44	MET
20	CT	57	ARG
20	CT	95	ALA
20	CT	97	ALA
21	CU	9	ARG
21	CU	12	LYS
25	CY	131	ASN
25	CY	134	ARG
25	CY	170	ALA
26	D0	2	ALA
28	D2	16	LEU
28	D2	36	ARG
29	D3	29	ARG
30	D4	44	THR
31	D5	14	ALA
31	D5	58	LEU
33	D7	32	LYS
34	D8	13	ARG
34	D8	36	LYS
34	D8	64	TYR
37	DC	24	GLU
37	DC	51	PRO
37	DC	146	GLY
37	DC	153	ILE
37	DC	172	HIS
37	DC	183	GLU
37	DC	201	PRO
37	DC	211	SER
38	DD	89	SER
38	DD	115	GLN
38	DD	157	ARG
38	DD	214	TRP
39	DE	43	GLY
39	DE	44	TYR
39	DE	60	ASN
39	DE	113	PHE
39	DE	162	ALA
39	DE	180	ASN
40	DF	14	PRO

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Mol	Chain	Res	Type
40	DF	33	LEU
40	DF	61	GLY
40	DF	90	PHE
40	DF	102	PRO
40	DF	115	ALA
40	DF	132	VAL
40	DF	133	ASN
41	DG	14	GLU
41	DG	175	LEU
42	DH	39	PRO
42	DH	134	SER
42	DH	152	ARG
42	DH	153	LYS
42	DH	154	PRO
43	DI	14	ASP
43	DI	40	THR
43	DI	85	GLU
43	DI	104	GLN
43	DI	115	ALA
43	DI	120	ILE
44	DN	18	ALA
44	DN	46	VAL
44	DN	57	ALA
44	DN	80	GLY
44	DN	129	PRO
45	DO	27	GLY
45	DO	36	GLY
45	DO	45	GLU
45	DO	55	GLY
45	DO	68	GLU
45	DO	90	GLN
46	DP	12	ALA
46	DP	110	TYR
46	DP	120	ALA
46	DP	129	ALA
46	DP	141	ALA
47	DQ	24	GLY
47	DQ	52	VAL
47	DQ	59	ARG
47	DQ	60	ARG
47	DQ	78	PRO
47	DQ	105	GLU

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Mol	Chain	Res	Type
47	DQ	115	MET
47	DQ	140	ALA
48	DR	5	LYS
48	DR	13	HIS
48	DR	17	ARG
48	DR	50	HIS
48	DR	60	LEU
48	DR	63	ARG
48	DR	88	ARG
48	DR	105	ARG
48	DR	107	ASP
49	DS	43	GLU
49	DS	53	SER
49	DS	54	LEU
49	DS	61	ASN
49	DS	73	LEU
49	DS	78	LEU
49	DS	85	VAL
50	DT	4	GLY
50	DT	25	GLY
50	DT	129	ARG
50	DT	131	ALA
51	DU	52	ARG
51	DU	83	LEU
51	DU	92	ARG
51	DU	99	ALA
52	DV	40	LEU
52	DV	42	GLY
52	DV	43	GLU
52	DV	48	GLY
52	DV	53	GLU
52	DV	90	PRO
53	DW	14	PRO
53	DW	66	GLU
53	DW	67	ASP
54	DX	7	VAL
54	DX	72	LYS
54	DX	89	ILE
55	DY	8	LYS
55	DY	23	ARG
55	DY	29	GLU
55	DY	31	LEU

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Mol	Chain	Res	Type
55	DY	40	GLU
55	DY	99	CYS
56	DZ	31	ARG
56	DZ	46	LYS
56	DZ	62	PRO
56	DZ	77	ASP
56	DZ	165	VAL
2	AB	97	TRP
2	AB	175	ARG
2	AB	217	ARG
3	AC	41	GLY
3	AC	179	ARG
4	AD	129	ASN
4	AD	142	PRO
5	AE	21	ALA
5	AE	70	PRO
5	AE	147	ASP
5	AE	154	GLY
6	AF	12	PRO
6	AF	22	GLU
6	AF	43	LEU
7	AG	58	PRO
7	AG	66	VAL
7	AG	104	LEU
7	AG	111	ARG
7	AG	119	ARG
7	AG	131	LYS
8	AH	54	ASP
9	AI	61	ALA
9	AI	89	ASN
9	AI	106	ALA
10	AJ	51	ARG
11	AK	18	ARG
11	AK	66	LEU
11	AK	72	ALA
12	AL	105	TYR
13	AM	46	LYS
15	AO	29	VAL
16	AP	7	ALA
16	AP	29	ASP
16	AP	61	SER
16	AP	65	GLN

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Mol	Chain	Res	Type
17	AQ	17	LYS
17	AQ	98	LEU
20	AT	24	LEU
20	AT	57	ARG
20	AT	65	LYS
25	AY	19	GLU
25	AY	171	LYS
25	AY	175	LEU
27	B1	12	PRO
27	B1	13	ILE
27	B1	85	LEU
28	B2	28	LYS
28	B2	43	GLN
30	B4	18	CYS
30	B4	40	HIS
31	B5	14	ALA
32	B6	23	THR
32	B6	28	ARG
33	B7	11	LYS
33	B7	33	ARG
34	B8	17	THR
34	B8	42	ARG
37	BC	73	ARG
37	BC	92	ASP
37	BC	151	GLU
37	BC	175	VAL
37	BC	185	LEU
37	BC	218	MET
38	BD	140	THR
38	BD	147	LEU
38	BD	211	ARG
39	BE	17	ASP
39	BE	64	LYS
39	BE	130	GLY
40	BF	38	ARG
40	BF	84	VAL
40	BF	90	PHE
40	BF	102	PRO
40	BF	146	ALA
41	BG	26	GLN
41	BG	107	LEU
41	BG	173	LEU

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Mol	Chain	Res	Type
41	BG	177	GLY
42	BH	15	VAL
42	BH	81	GLU
42	BH	98	LEU
42	BH	134	SER
42	BH	150	ALA
43	BI	14	ASP
43	BI	42	SER
43	BI	70	GLU
43	BI	75	LEU
43	BI	115	ALA
43	BI	132	PRO
44	BN	15	LEU
44	BN	56	ASN
44	BN	94	HIS
45	BO	5	GLN
45	BO	55	GLY
45	BO	108	GLU
45	BO	110	GLY
45	BO	120	GLU
46	BP	109	GLY
47	BQ	13	GLN
47	BQ	54	MET
48	BR	13	HIS
48	BR	63	ARG
48	BR	81	ASP
48	BR	102	GLU
49	BS	66	ALA
49	BS	88	ASP
50	BT	4	GLY
50	BT	55	ASN
50	BT	91	ARG
50	BT	95	ARG
50	BT	129	ARG
51	BU	52	ARG
51	BU	90	VAL
51	BU	99	ALA
52	BV	32	THR
52	BV	90	PRO
53	BW	2	GLU
53	BW	67	ASP
54	BX	48	LYS

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Mol	Chain	Res	Type
54	BX	50	LYS
54	BX	72	LYS
55	BY	8	LYS
55	BY	11	ASP
55	BY	17	SER
55	BY	21	LYS
55	BY	87	LYS
56	BZ	49	ARG
56	BZ	56	VAL
56	BZ	108	PRO
56	BZ	178	GLU
2	CB	205	ASP
3	CC	41	GLY
3	CC	45	LYS
4	CD	40	PRO
4	CD	43	HIS
4	CD	142	PRO
4	CD	153	ARG
5	CE	21	ALA
5	CE	70	PRO
5	CE	147	ASP
5	CE	154	GLY
6	CF	12	PRO
6	CF	43	LEU
6	CF	86	ARG
7	CG	58	PRO
7	CG	66	VAL
7	CG	77	SER
7	CG	104	LEU
7	CG	111	ARG
7	CG	136	LYS
8	CH	54	ASP
8	CH	63	LEU
8	CH	68	ARG
9	CI	61	ALA
9	CI	89	ASN
9	CI	106	ALA
10	CJ	51	ARG
11	CK	72	ALA
12	CL	46	LYS
12	CL	105	TYR
13	CM	46	LYS

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Mol	Chain	Res	Type
13	CM	81	LEU
15	CO	37	ASN
15	CO	88	ARG
16	CP	7	ALA
16	CP	29	ASP
16	CP	65	GLN
16	CP	83	GLU
17	CQ	98	LEU
20	CT	12	ALA
20	CT	22	ARG
20	CT	24	LEU
20	CT	65	LYS
20	CT	96	GLY
25	CY	12	SER
25	CY	16	LYS
25	CY	17	SER
25	CY	18	LEU
25	CY	54	GLN
25	CY	84	ARG
25	CY	88	LEU
25	CY	108	GLU
25	CY	112	LYS
25	CY	121	TYR
25	CY	140	LEU
25	CY	158	GLU
25	CY	164	ILE
25	CY	172	ALA
27	D1	27	GLU
27	D1	70	VAL
30	D4	40	HIS
31	D5	20	ARG
32	D6	23	THR
32	D6	28	ARG
34	D8	17	THR
34	D8	47	LYS
34	D8	50	LEU
34	D8	51	ALA
37	DC	70	LYS
37	DC	92	ASP
37	DC	151	GLU
37	DC	175	VAL
37	DC	185	LEU

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Mol	Chain	Res	Type
37	DC	214	VAL
37	DC	218	MET
38	DD	140	THR
38	DD	211	ARG
38	DD	256	GLY
39	DE	17	ASP
39	DE	64	LYS
39	DE	130	GLY
40	DF	64	ILE
40	DF	84	VAL
40	DF	146	ALA
40	DF	164	ARG
40	DF	179	GLU
41	DG	35	GLU
41	DG	43	LEU
41	DG	45	GLU
41	DG	49	ASP
41	DG	117	PHE
42	DH	15	VAL
42	DH	81	GLU
42	DH	98	LEU
42	DH	148	ILE
42	DH	150	ALA
43	DI	6	LEU
43	DI	70	GLU
43	DI	75	LEU
43	DI	76	THR
43	DI	132	PRO
44	DN	15	LEU
44	DN	28	THR
44	DN	56	ASN
44	DN	94	HIS
45	DO	5	GLN
45	DO	108	GLU
45	DO	110	GLY
45	DO	114	ILE
46	DP	33	ARG
46	DP	52	GLU
46	DP	109	GLY
47	DQ	8	LYS
47	DQ	54	MET
48	DR	81	ASP

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Mol	Chain	Res	Type
48	DR	102	GLU
49	DS	66	ALA
49	DS	88	ASP
49	DS	97	ARG
50	DT	23	ARG
50	DT	55	ASN
50	DT	91	ARG
50	DT	117	ASP
51	DU	71	GLN
51	DU	88	ILE
51	DU	91	ASP
52	DV	32	THR
52	DV	82	ARG
53	DW	18	ARG
54	DX	15	GLU
54	DX	48	LYS
55	DY	11	ASP
55	DY	17	SER
55	DY	19	LYS
55	DY	26	LYS
55	DY	39	VAL
55	DY	87	LYS
56	DZ	51	ALA
56	DZ	138	GLU
56	DZ	139	VAL
56	DZ	140	ASP
2	AB	46	LYS
2	AB	131	PRO
2	AB	205	ASP
3	AC	9	GLY
3	AC	45	LYS
3	AC	109	PRO
4	AD	39	PRO
4	AD	40	PRO
4	AD	102	ASP
5	AE	128	PRO
6	AF	86	ARG
7	AG	23	VAL
7	AG	49	ILE
7	AG	77	SER
7	AG	96	GLN
7	AG	103	TRP

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Mol	Chain	Res	Type
7	AG	136	LYS
8	AH	87	SER
9	AI	57	GLY
11	AK	128	ALA
12	AL	9	GLN
13	AM	37	THR
13	AM	108	ARG
14	AN	7	ILE
15	AO	21	ASP
15	AO	38	ARG
15	AO	88	ARG
16	AP	2	VAL
16	AP	14	ASN
17	AQ	68	ARG
18	AR	72	ARG
19	AS	73	GLU
20	AT	22	ARG
20	AT	33	ILE
20	AT	34	LYS
20	AT	56	MET
25	AY	32	ARG
25	AY	49	HIS
25	AY	82	ALA
25	AY	104	PRO
26	B0	57	PHE
26	B0	83	PRO
28	B2	61	LEU
33	B7	35	ARG
34	B8	50	LEU
37	BC	55	ASP
37	BC	64	LEU
37	BC	70	LYS
37	BC	169	GLY
37	BC	209	LEU
38	BD	157	ARG
38	BD	256	GLY
39	BE	113	PHE
39	BE	149	ARG
40	BF	164	ARG
40	BF	177	ALA
41	BG	10	LYS
41	BG	56	ALA

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Mol	Chain	Res	Type
41	BG	69	ALA
41	BG	105	LYS
42	BH	110	SER
43	BI	76	THR
44	BN	28	THR
45	BO	79	PHE
46	BP	52	GLU
46	BP	59	LEU
46	BP	71	VAL
46	BP	115	LEU
47	BQ	8	LYS
47	BQ	50	ALA
48	BR	32	GLY
48	BR	39	PRO
48	BR	40	LYS
49	BS	19	LYS
49	BS	57	LYS
50	BT	117	ASP
51	BU	71	GLN
51	BU	111	GLU
52	BV	58	VAL
52	BV	74	LYS
53	BW	15	ARG
53	BW	44	ALA
53	BW	45	TYR
53	BW	110	LYS
54	BX	15	GLU
54	BX	38	GLU
54	BX	85	PRO
55	BY	50	ARG
55	BY	77	PRO
56	BZ	5	LEU
56	BZ	18	LEU
2	CB	25	ASN
2	CB	97	TRP
2	CB	131	PRO
2	CB	175	ARG
2	CB	217	ARG
3	CC	9	GLY
3	CC	109	PRO
5	CE	128	PRO
6	CF	22	GLU

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Mol	Chain	Res	Type
7	CG	23	VAL
7	CG	49	ILE
7	CG	103	TRP
9	CI	57	GLY
11	CK	128	ALA
12	CL	61	THR
13	CM	47	ASP
13	CM	108	ARG
15	CO	21	ASP
15	CO	29	VAL
15	CO	38	ARG
15	CO	64	ARG
16	CP	2	VAL
16	CP	61	SER
18	CR	27	GLY
19	CS	73	GLU
20	CT	23	ARG
20	CT	33	ILE
20	CT	56	MET
20	CT	79	ARG
25	CY	7	TYR
25	CY	122	ALA
25	CY	159	ALA
25	CY	162	GLN
25	CY	163	LYS
26	D0	57	PHE
26	D0	83	PRO
27	D1	58	ILE
27	D1	59	THR
30	D4	18	CYS
31	D5	3	LYS
33	D7	7	PRO
33	D7	11	LYS
33	D7	19	ARG
37	DC	55	ASP
37	DC	64	LEU
37	DC	73	ARG
37	DC	169	GLY
37	DC	209	LEU
38	DD	147	LEU
38	DD	170	GLY
39	DE	69	LYS

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Mol	Chain	Res	Type
39	DE	72	VAL
39	DE	149	ARG
40	DF	177	ALA
41	DG	101	ILE
41	DG	133	LEU
41	DG	138	GLN
43	DI	42	SER
43	DI	87	LYS
43	DI	119	PRO
44	DN	40	PRO
45	DO	43	VAL
45	DO	79	PHE
46	DP	59	LEU
46	DP	76	LYS
46	DP	115	LEU
47	DQ	21	THR
47	DQ	50	ALA
48	DR	32	GLY
49	DS	19	LYS
49	DS	57	LYS
50	DT	95	ARG
51	DU	54	LYS
51	DU	61	TRP
51	DU	90	VAL
52	DV	58	VAL
52	DV	74	LYS
53	DW	2	GLU
53	DW	65	LEU
54	DX	50	LYS
55	DY	50	ARG
55	DY	77	PRO
56	DZ	22	GLY
56	DZ	32	HIS
56	DZ	63	ASP
56	DZ	104	PHE
56	DZ	177	PRO
2	AB	25	ASN
3	AC	8	ILE
3	AC	15	THR
3	AC	81	GLY
5	AE	77	PRO
8	AH	49	GLU

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Mol	Chain	Res	Type
8	AH	76	PRO
9	AI	34	ASN
13	AM	6	GLY
13	AM	32	GLU
13	AM	116	THR
14	AN	56	VAL
15	AO	64	ARG
18	AR	27	GLY
20	AT	23	ARG
20	AT	79	ARG
25	AY	70	SER
31	B5	28	PRO
31	B5	38	ALA
38	BD	170	GLY
38	BD	242	ARG
39	BE	69	LYS
39	BE	72	VAL
39	BE	98	PRO
40	BF	83	PHE
41	BG	15	VAL
41	BG	142	PRO
42	BH	93	GLY
43	BI	87	LYS
43	BI	119	PRO
44	BN	40	PRO
44	BN	125	GLY
46	BP	26	GLY
46	BP	76	LYS
46	BP	106	LEU
47	BQ	10	ARG
47	BQ	52	VAL
47	BQ	56	ARG
48	BR	73	VAL
49	BS	103	GLU
50	BT	82	LEU
51	BU	54	LYS
51	BU	61	TRP
52	BV	3	ALA
52	BV	9	GLY
53	BW	65	LEU
55	BY	26	LYS
55	BY	81	LYS

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Mol	Chain	Res	Type
56	BZ	37	VAL
56	BZ	106	GLY
2	CB	177	ALA
2	CB	211	ILE
3	CC	8	ILE
3	CC	15	THR
3	CC	43	LEU
3	CC	81	GLY
3	CC	173	VAL
4	CD	39	PRO
5	CE	76	ILE
5	CE	77	PRO
7	CG	96	GLN
8	CH	30	ARG
8	CH	76	PRO
8	CH	87	SER
9	CI	34	ASN
9	CI	126	SER
12	CL	9	GLN
12	CL	13	LYS
12	CL	101	VAL
13	CM	6	GLY
13	CM	37	THR
14	CN	56	VAL
16	CP	14	ASN
17	CQ	68	ARG
25	CY	31	GLY
25	CY	92	PRO
30	D4	36	CYS
31	D5	24	ALA
31	D5	28	PRO
32	D6	51	GLU
39	DE	134	ILE
40	DF	38	ARG
40	DF	83	PHE
41	DG	24	GLY
41	DG	74	LYS
41	DG	171	ALA
42	DH	93	GLY
44	DN	125	GLY
45	DO	63	VAL
46	DP	71	VAL

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Mol	Chain	Res	Type
46	DP	106	LEU
47	DQ	13	GLN
48	DR	39	PRO
48	DR	73	VAL
51	DU	112	ARG
54	DX	38	GLU
54	DX	85	PRO
55	DY	21	LYS
55	DY	81	LYS
2	AB	130	ARG
3	AC	55	VAL
3	AC	173	VAL
4	AD	178	VAL
9	AI	98	PRO
11	AK	49	GLY
12	AL	101	VAL
27	B1	79	GLY
30	B4	19	GLY
34	B8	20	GLY
37	BC	143	GLY
38	BD	118	VAL
38	BD	244	ARG
38	BD	246	PRO
42	BH	127	GLU
43	BI	84	GLY
45	BO	63	VAL
48	BR	106	GLY
52	BV	99	ILE
56	BZ	82	ARG
2	CB	130	ARG
3	CC	55	VAL
4	CD	95	GLY
4	CD	178	VAL
9	CI	98	PRO
11	CK	49	GLY
30	D4	19	GLY
34	D8	20	GLY
37	DC	143	GLY
37	DC	200	LYS
38	DD	10	THR
38	DD	118	VAL
38	DD	244	ARG

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Mol	Chain	Res	Type
39	DE	34	VAL
39	DE	98	PRO
40	DF	171	PRO
41	DG	39	ILE
41	DG	67	LYS
46	DP	26	GLY
52	DV	9	GLY
52	DV	99	ILE
56	DZ	116	VAL
2	AB	211	ILE
3	AC	108	ASN
7	AG	88	PRO
27	B1	31	GLY
31	B5	47	PRO
33	B7	7	PRO
37	BC	19	VAL
37	BC	181	PRO
37	BC	200	LYS
38	BD	10	THR
39	BE	34	VAL
39	BE	134	ILE
45	BO	43	VAL
49	BS	60	GLY
55	BY	37	VAL
56	BZ	22	GLY
3	CC	108	ASN
4	CD	56	VAL
6	CF	67	MET
9	CI	108	VAL
11	CK	115	PRO
14	CN	7	ILE
37	DC	19	VAL
37	DC	181	PRO
38	DD	246	PRO
42	DH	29	PRO
42	DH	127	GLU
43	DI	84	GLY
48	DR	106	GLY
49	DS	60	GLY
53	DW	26	GLY
55	DY	37	VAL
4	AD	148	VAL

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Mol	Chain	Res	Type
7	AG	17	VAL
8	AH	128	GLY
9	AI	108	VAL
10	AJ	24	VAL
25	AY	35	PRO
27	B1	68	PRO
37	BC	132	ALA
37	BC	180	PHE
39	BE	39	PRO
40	BF	171	PRO
41	BG	77	ILE
42	BH	29	PRO
46	BP	43	GLY
48	BR	85	PRO
49	BS	28	VAL
53	BW	26	GLY
7	CG	17	VAL
7	CG	88	PRO
10	CJ	24	VAL
27	D1	28	GLY
30	D4	15	ILE
31	D5	47	PRO
37	DC	132	ALA
37	DC	180	PHE
39	DE	39	PRO
41	DG	149	VAL
46	DP	97	PRO
49	DS	28	VAL
49	DS	104	GLY
54	DX	31	HIS
56	DZ	82	ARG
56	DZ	167	PRO
4	AD	8	VAL
5	AE	76	ILE
6	AF	67	MET
8	AH	51	VAL
11	AK	115	PRO
15	AO	86	GLY
27	B1	70	VAL
30	B4	15	ILE
32	B6	41	PRO
38	BD	152	GLY

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Mol	Chain	Res	Type
41	BG	92	VAL
42	BH	22	GLY
43	BI	23	PRO
43	BI	127	VAL
45	BO	101	PRO
46	BP	37	GLY
46	BP	97	PRO
49	BS	104	GLY
52	BV	70	ILE
53	BW	59	VAL
54	BX	31	HIS
55	BY	53	PRO
56	BZ	68	PRO
2	CB	230	VAL
4	CD	148	VAL
8	CH	51	VAL
8	CH	128	GLY
10	CJ	93	GLY
25	CY	61	PRO
28	D2	42	GLY
33	D7	17	GLY
39	DE	116	VAL
40	DF	81	PRO
41	DG	89	GLY
45	DO	101	PRO
46	DP	43	GLY
47	DQ	127	ILE
48	DR	85	PRO
52	DV	70	ILE
53	DW	59	VAL
55	DY	10	GLY
55	DY	53	PRO
48	BR	46	GLY
50	BT	88	ILE
53	BW	20	VAL
4	CD	8	VAL
38	DD	152	GLY
43	DI	127	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	AB	202/220 (92%)	177 (88%)	25 (12%)	4	23
2	CB	202/220 (92%)	177 (88%)	25 (12%)	4	23
3	AC	160/188 (85%)	142 (89%)	18 (11%)	6	27
3	CC	160/188 (85%)	142 (89%)	18 (11%)	6	27
4	AD	180/181 (99%)	155 (86%)	25 (14%)	3	20
4	CD	180/181 (99%)	156 (87%)	24 (13%)	4	21
5	AE	115/123 (94%)	99 (86%)	16 (14%)	3	20
5	CE	115/123 (94%)	99 (86%)	16 (14%)	3	20
6	AF	90/90 (100%)	86 (96%)	4 (4%)	28	62
6	CF	90/90 (100%)	86 (96%)	4 (4%)	28	62
7	AG	126/127 (99%)	118 (94%)	8 (6%)	18	51
7	CG	126/127 (99%)	116 (92%)	10 (8%)	12	41
8	AH	119/119 (100%)	108 (91%)	11 (9%)	9	36
8	CH	119/119 (100%)	108 (91%)	11 (9%)	9	36
9	AI	98/99 (99%)	83 (85%)	15 (15%)	2	17
9	CI	98/99 (99%)	85 (87%)	13 (13%)	4	21
10	AJ	88/92 (96%)	81 (92%)	7 (8%)	12	41
10	CJ	88/92 (96%)	81 (92%)	7 (8%)	12	41
11	AK	90/99 (91%)	80 (89%)	10 (11%)	6	28
11	CK	90/99 (91%)	81 (90%)	9 (10%)	7	32
12	AL	104/111 (94%)	91 (88%)	13 (12%)	4	23
12	CL	104/111 (94%)	93 (89%)	11 (11%)	6	30
13	AM	99/101 (98%)	92 (93%)	7 (7%)	14	46
13	CM	99/101 (98%)	92 (93%)	7 (7%)	14	46
14	AN	49/50 (98%)	45 (92%)	4 (8%)	11	40
14	CN	49/50 (98%)	45 (92%)	4 (8%)	11	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
15	AO	79/80 (99%)	71 (90%)	8 (10%)	7	32
15	CO	79/80 (99%)	70 (89%)	9 (11%)	5	26
16	AP	72/74 (97%)	61 (85%)	11 (15%)	2	17
16	CP	72/74 (97%)	61 (85%)	11 (15%)	2	17
17	AQ	94/97 (97%)	88 (94%)	6 (6%)	17	50
17	CQ	94/97 (97%)	88 (94%)	6 (6%)	17	50
18	AR	61/77 (79%)	56 (92%)	5 (8%)	11	40
18	CR	61/77 (79%)	55 (90%)	6 (10%)	8	33
19	AS	69/80 (86%)	57 (83%)	12 (17%)	2	11
19	CS	69/80 (86%)	56 (81%)	13 (19%)	1	8
20	AT	76/82 (93%)	65 (86%)	11 (14%)	3	18
20	CT	76/82 (93%)	65 (86%)	11 (14%)	3	18
21	AU	19/22 (86%)	18 (95%)	1 (5%)	22	55
21	CU	19/22 (86%)	18 (95%)	1 (5%)	22	55
25	AY	157/157 (100%)	135 (86%)	22 (14%)	3	19
25	CY	157/157 (100%)	142 (90%)	15 (10%)	8	34
26	B0	61/67 (91%)	57 (93%)	4 (7%)	16	49
26	D0	61/67 (91%)	57 (93%)	4 (7%)	16	49
27	B1	73/83 (88%)	54 (74%)	19 (26%)	0	3
27	D1	73/83 (88%)	50 (68%)	23 (32%)	0	2
28	B2	46/67 (69%)	31 (67%)	15 (33%)	0	2
28	D2	46/67 (69%)	34 (74%)	12 (26%)	0	3
29	B3	51/52 (98%)	48 (94%)	3 (6%)	19	53
29	D3	51/52 (98%)	48 (94%)	3 (6%)	19	53
31	B5	51/52 (98%)	42 (82%)	9 (18%)	2	10
31	D5	51/52 (98%)	43 (84%)	8 (16%)	2	15
32	B6	43/52 (83%)	34 (79%)	9 (21%)	1	6
32	D6	43/52 (83%)	34 (79%)	9 (21%)	1	6
33	B7	41/42 (98%)	34 (83%)	7 (17%)	2	12
33	D7	41/42 (98%)	34 (83%)	7 (17%)	2	12
34	B8	53/55 (96%)	45 (85%)	8 (15%)	3	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	D8	53/55 (96%)	45 (85%)	8 (15%)	3	17
37	BC	61/181 (34%)	56 (92%)	5 (8%)	11	40
37	DC	61/181 (34%)	57 (93%)	4 (7%)	16	49
38	BD	213/218 (98%)	186 (87%)	27 (13%)	4	22
38	DD	213/218 (98%)	186 (87%)	27 (13%)	4	22
39	BE	165/166 (99%)	142 (86%)	23 (14%)	3	20
39	DE	165/166 (99%)	141 (86%)	24 (14%)	3	18
40	BF	165/166 (99%)	142 (86%)	23 (14%)	3	20
40	DF	165/166 (99%)	142 (86%)	23 (14%)	3	20
41	BG	155/156 (99%)	137 (88%)	18 (12%)	5	26
41	DG	155/156 (99%)	127 (82%)	28 (18%)	1	9
42	BH	132/148 (89%)	115 (87%)	17 (13%)	4	22
42	DH	132/148 (89%)	116 (88%)	16 (12%)	5	24
43	BI	122/124 (98%)	103 (84%)	19 (16%)	2	16
43	DI	122/124 (98%)	103 (84%)	19 (16%)	2	16
44	BN	117/119 (98%)	100 (86%)	17 (14%)	3	18
44	DN	117/119 (98%)	99 (85%)	18 (15%)	2	16
45	BO	100/100 (100%)	80 (80%)	20 (20%)	1	7
45	DO	100/100 (100%)	79 (79%)	21 (21%)	1	5
46	BP	112/116 (97%)	84 (75%)	28 (25%)	0	4
46	DP	112/116 (97%)	84 (75%)	28 (25%)	0	4
47	BQ	106/111 (96%)	86 (81%)	20 (19%)	1	8
47	DQ	106/111 (96%)	88 (83%)	18 (17%)	2	12
48	BR	100/101 (99%)	80 (80%)	20 (20%)	1	7
48	DR	100/101 (99%)	80 (80%)	20 (20%)	1	7
49	BS	77/88 (88%)	63 (82%)	14 (18%)	1	9
49	DS	77/88 (88%)	62 (80%)	15 (20%)	1	7
50	BT	120/127 (94%)	89 (74%)	31 (26%)	0	3
50	DT	120/127 (94%)	89 (74%)	31 (26%)	0	3
51	BU	92/94 (98%)	80 (87%)	12 (13%)	4	21
51	DU	92/94 (98%)	81 (88%)	11 (12%)	5	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
52	BV	82/82 (100%)	68 (83%)	14 (17%)	2	12
52	DV	82/82 (100%)	68 (83%)	14 (17%)	2	12
53	BW	91/92 (99%)	78 (86%)	13 (14%)	3	19
53	DW	91/92 (99%)	78 (86%)	13 (14%)	3	19
54	BX	74/78 (95%)	59 (80%)	15 (20%)	1	6
54	DX	74/78 (95%)	58 (78%)	16 (22%)	1	5
55	BY	84/91 (92%)	63 (75%)	21 (25%)	0	4
55	DY	84/91 (92%)	63 (75%)	21 (25%)	0	4
56	BZ	155/179 (87%)	134 (86%)	21 (14%)	4	21
56	DZ	155/179 (87%)	124 (80%)	31 (20%)	1	7
All	All	9778/10552 (93%)	8384 (86%)	1394 (14%)	3	19

All (1394) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	AB	17	PHE
2	AB	22	LYS
2	AB	31	TYR
2	AB	33	TYR
2	AB	44	LEU
2	AB	56	ARG
2	AB	67	THR
2	AB	69	LEU
2	AB	70	PHE
2	AB	75	LYS
2	AB	80	ILE
2	AB	96	ARG
2	AB	97	TRP
2	AB	98	LEU
2	AB	102	LEU
2	AB	108	ILE
2	AB	115	LEU
2	AB	116	GLU
2	AB	137	ARG
2	AB	140	HIS
2	AB	158	LEU
2	AB	178	ARG
2	AB	187	LEU

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Mol	Chain	Res	Type
2	AB	212	GLN
2	AB	222	ILE
3	AC	5	ILE
3	AC	12	LEU
3	AC	16	ARG
3	AC	27	LYS
3	AC	37	GLN
3	AC	43	LEU
3	AC	52	LEU
3	AC	101	LEU
3	AC	104	GLN
3	AC	120	VAL
3	AC	127	ARG
3	AC	131	ARG
3	AC	156	ARG
3	AC	164	ARG
3	AC	165	THR
3	AC	166	GLU
3	AC	167	TRP
3	AC	181	ASN
4	AD	3	ARG
4	AD	9	CYS
4	AD	21	LEU
4	AD	26	CYS
4	AD	31	CYS
4	AD	33	MET
4	AD	42	GLN
4	AD	53	ASP
4	AD	59	ARG
4	AD	79	PHE
4	AD	96	LEU
4	AD	97	LEU
4	AD	110	PHE
4	AD	120	LEU
4	AD	122	ARG
4	AD	131	ARG
4	AD	132	ARG
4	AD	138	TYR
4	AD	158	ILE
4	AD	170	VAL
4	AD	181	MET
4	AD	194	LEU

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Mol	Chain	Res	Type
4	AD	196	LEU
4	AD	200	GLU
4	AD	202	LEU
5	AE	10	MET
5	AE	12	LEU
5	AE	16	THR
5	AE	20	GLN
5	AE	47	LYS
5	AE	60	TYR
5	AE	61	TYR
5	AE	68	GLU
5	AE	71	LEU
5	AE	73	ASN
5	AE	76	ILE
5	AE	90	VAL
5	AE	96	PRO
5	AE	101	ILE
5	AE	131	ILE
5	AE	149	GLU
6	AF	32	ASN
6	AF	75	LEU
6	AF	98	LEU
6	AF	100	ASN
7	AG	14	PRO
7	AG	41	ARG
7	AG	54	THR
7	AG	57	GLU
7	AG	84	ASN
7	AG	111	ARG
7	AG	151	TYR
7	AG	156	TRP
8	AH	1	MET
8	AH	3	THR
8	AH	10	LEU
8	AH	18	ARG
8	AH	25	ASP
8	AH	35	ILE
8	AH	102	ARG
8	AH	119	LEU
8	AH	133	LEU
8	AH	136	GLU
8	AH	137	VAL

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Mol	Chain	Res	Type
9	AI	4	TYR
9	AI	27	THR
9	AI	63	ILE
9	AI	66	ARG
9	AI	78	LYS
9	AI	89	ASN
9	AI	95	LYS
9	AI	101	PHE
9	AI	104	ARG
9	AI	108	VAL
9	AI	113	LYS
9	AI	114	TYR
9	AI	121	ARG
9	AI	125	TYR
9	AI	128	ARG
10	AJ	16	LEU
10	AJ	22	LYS
10	AJ	47	PHE
10	AJ	50	ILE
10	AJ	62	HIS
10	AJ	84	GLN
10	AJ	96	ILE
11	AK	18	ARG
11	AK	21	ILE
11	AK	48	ILE
11	AK	91	ARG
11	AK	103	LEU
11	AK	105	VAL
11	AK	117	ASN
11	AK	125	PHE
11	AK	126	ARG
11	AK	127	LYS
12	AL	19	ARG
12	AL	20	LYS
12	AL	21	LYS
12	AL	41	ARG
12	AL	48	PRO
12	AL	49	ASN
12	AL	53	ARG
12	AL	66	VAL
12	AL	70	ILE
12	AL	79	GLU

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Mol	Chain	Res	Type
12	AL	84	LEU
12	AL	89	ARG
12	AL	102	ARG
13	AM	47	ASP
13	AM	48	LEU
13	AM	64	TRP
13	AM	82	MET
13	AM	93	ARG
13	AM	108	ARG
13	AM	120	LYS
14	AN	16	PHE
14	AN	26	ARG
14	AN	29	ARG
14	AN	61	TRP
15	AO	4	THR
15	AO	6	GLU
15	AO	10	LYS
15	AO	35	ARG
15	AO	46	HIS
15	AO	47	LYS
15	AO	65	ARG
15	AO	82	ILE
16	AP	1	MET
16	AP	6	LEU
16	AP	26	ARG
16	AP	27	LYS
16	AP	29	ASP
16	AP	38	TYR
16	AP	39	TYR
16	AP	48	TRP
16	AP	69	THR
16	AP	73	LEU
16	AP	82	GLN
17	AQ	36	ILE
17	AQ	55	ASP
17	AQ	56	VAL
17	AQ	78	GLU
17	AQ	81	ARG
17	AQ	89	LEU
18	AR	29	PHE
18	AR	31	LEU
18	AR	46	GLU

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Mol	Chain	Res	Type
18	AR	58	LEU
18	AR	74	ARG
19	AS	5	LEU
19	AS	6	LYS
19	AS	13	ASP
19	AS	27	GLU
19	AS	34	TRP
19	AS	37	ARG
19	AS	43	GLU
19	AS	44	MET
19	AS	49	ILE
19	AS	53	ASN
19	AS	70	LYS
19	AS	78	ARG
20	AT	8	ARG
20	AT	13	LEU
20	AT	24	LEU
20	AT	26	ASN
20	AT	57	ARG
20	AT	62	LEU
20	AT	71	THR
20	AT	73	HIS
20	AT	74	LYS
20	AT	93	GLU
20	AT	100	ILE
21	AU	24	ARG
25	AY	13	HIS
25	AY	20	VAL
25	AY	29	ARG
25	AY	41	LEU
25	AY	44	GLU
25	AY	53	ASN
25	AY	55	ILE
25	AY	64	ARG
25	AY	76	LEU
25	AY	91	ASN
25	AY	101	ILE
25	AY	103	ILE
25	AY	113	ASP
25	AY	130	ARG
25	AY	138	ASP
25	AY	142	LYS

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Mol	Chain	Res	Type
25	AY	150	SER
25	AY	152	ASP
25	AY	164	ILE
25	AY	168	PHE
25	AY	169	ILE
25	AY	174	GLN
26	B0	20	ARG
26	B0	38	VAL
26	B0	44	ARG
26	B0	64	ASP
27	B1	19	GLN
27	B1	20	ARG
27	B1	21	ARG
27	B1	26	ARG
27	B1	34	THR
27	B1	39	LYS
27	B1	46	LEU
27	B1	47	GLN
27	B1	48	LYS
27	B1	53	VAL
27	B1	56	GLN
27	B1	58	ILE
27	B1	59	THR
27	B1	62	VAL
27	B1	69	LYS
27	B1	78	LYS
27	B1	87	PRO
27	B1	89	GLU
27	B1	94	LEU
28	B2	14	ARG
28	B2	16	LEU
28	B2	17	SER
28	B2	22	GLU
28	B2	26	ARG
28	B2	27	GLU
28	B2	30	ARG
28	B2	32	LEU
28	B2	33	MET
28	B2	44	LEU
28	B2	46	GLN
28	B2	47	ASN
28	B2	51	ARG

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Mol	Chain	Res	Type
28	B2	56	GLN
28	B2	57	ILE
29	B3	8	LEU
29	B3	26	LEU
29	B3	31	LEU
31	B5	3	LYS
31	B5	4	HIS
31	B5	5	PRO
31	B5	22	HIS
31	B5	32	PRO
31	B5	43	HIS
31	B5	48	GLU
31	B5	56	LYS
31	B5	58	LEU
32	B6	10	LEU
32	B6	20	ASN
32	B6	31	PRO
32	B6	35	GLU
32	B6	37	ARG
32	B6	42	TRP
32	B6	43	CYS
32	B6	44	ARG
32	B6	51	GLU
33	B7	8	ASN
33	B7	24	THR
33	B7	36	GLN
33	B7	43	THR
33	B7	44	PRO
33	B7	46	VAL
33	B7	48	LYS
34	B8	31	HIS
34	B8	40	GLU
34	B8	44	LYS
34	B8	46	ARG
34	B8	48	PHE
34	B8	52	LYS
34	B8	60	LEU
34	B8	64	TYR
37	BC	36	LYS
37	BC	37	PHE
37	BC	49	ILE
37	BC	64	LEU

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Mol	Chain	Res	Type
37	BC	66	HIS
38	BD	5	LYS
38	BD	10	THR
38	BD	14	ARG
38	BD	15	PHE
38	BD	24	ILE
38	BD	26	LYS
38	BD	31	LYS
38	BD	46	GLN
38	BD	71	ASP
38	BD	91	ARG
38	BD	94	LEU
38	BD	95	LEU
38	BD	96	HIS
38	BD	131	LEU
38	BD	135	PHE
38	BD	166	GLN
38	BD	173	VAL
38	BD	189	CYS
38	BD	190	TYR
38	BD	198	ASN
38	BD	211	ARG
38	BD	221	VAL
38	BD	228	PRO
38	BD	241	PRO
38	BD	255	LYS
38	BD	261	LYS
38	BD	271	ILE
39	BE	21	VAL
39	BE	44	TYR
39	BE	61	ARG
39	BE	64	LYS
39	BE	67	PHE
39	BE	69	LYS
39	BE	78	LEU
39	BE	79	ARG
39	BE	113	PHE
39	BE	118	LYS
39	BE	119	ARG
39	BE	122	PHE
39	BE	127	ASP
39	BE	134	ILE

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Mol	Chain	Res	Type
39	BE	141	ILE
39	BE	143	ASN
39	BE	178	GLU
39	BE	181	LEU
39	BE	183	LEU
39	BE	184	VAL
39	BE	192	ASN
39	BE	202	LYS
39	BE	203	LYS
40	BF	6	VAL
40	BF	23	ASP
40	BF	24	LEU
40	BF	28	ILE
40	BF	46	ARG
40	BF	50	SER
40	BF	53	THR
40	BF	56	GLU
40	BF	65	TRP
40	BF	66	PRO
40	BF	67	GLN
40	BF	70	THR
40	BF	74	ARG
40	BF	77	ASP
40	BF	104	LYS
40	BF	112	MET
40	BF	116	ASP
40	BF	136	THR
40	BF	164	ARG
40	BF	169	ASN
40	BF	197	ASP
40	BF	201	VAL
40	BF	202	PHE
41	BG	22	ARG
41	BG	33	ARG
41	BG	43	LEU
41	BG	45	GLU
41	BG	51	ARG
41	BG	60	LEU
41	BG	64	THR
41	BG	77	ILE
41	BG	87	PRO
41	BG	106	LEU

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Mol	Chain	Res	Type
41	BG	125	PHE
41	BG	133	LEU
41	BG	136	ARG
41	BG	139	LEU
41	BG	147	ASP
41	BG	159	VAL
41	BG	170	ARG
41	BG	174	GLU
42	BH	41	MET
42	BH	53	GLU
42	BH	61	HIS
42	BH	68	THR
42	BH	83	TYR
42	BH	86	GLU
42	BH	89	ILE
42	BH	105	LEU
42	BH	123	PHE
42	BH	136	ILE
42	BH	143	GLN
42	BH	153	LYS
42	BH	157	TYR
42	BH	158	HIS
42	BH	159	GLU
42	BH	163	TYR
42	BH	170	ARG
43	BI	1	MET
43	BI	3	VAL
43	BI	5	LEU
43	BI	7	GLU
43	BI	10	GLU
43	BI	12	LEU
43	BI	25	TYR
43	BI	38	LEU
43	BI	51	ILE
43	BI	54	GLN
43	BI	64	GLU
43	BI	86	THR
43	BI	96	ASP
43	BI	118	LYS
43	BI	122	GLU
43	BI	123	LEU
43	BI	126	TYR

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Mol	Chain	Res	Type
43	BI	130	TYR
43	BI	136	VAL
44	BN	4	TYR
44	BN	19	GLU
44	BN	34	LEU
44	BN	39	ARG
44	BN	42	TRP
44	BN	56	ASN
44	BN	62	VAL
44	BN	66	LYS
44	BN	82	LEU
44	BN	87	LEU
44	BN	94	HIS
44	BN	99	LEU
44	BN	109	LYS
44	BN	123	TYR
44	BN	127	ASP
44	BN	128	HIS
44	BN	131	GLN
45	BO	1	MET
45	BO	7	TYR
45	BO	9	GLU
45	BO	19	ILE
45	BO	24	VAL
45	BO	25	LEU
45	BO	32	TYR
45	BO	39	ILE
45	BO	45	GLU
45	BO	47	ILE
45	BO	48	PRO
45	BO	69	ILE
45	BO	73	ASP
45	BO	80	ASP
45	BO	82	ASN
45	BO	86	ILE
45	BO	87	ILE
45	BO	91	LEU
45	BO	102	VAL
45	BO	115	VAL
46	BP	6	LEU
46	BP	13	ASN
46	BP	16	ARG

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Mol	Chain	Res	Type
46	BP	18	ARG
46	BP	21	ARG
46	BP	32	THR
46	BP	35	HIS
46	BP	50	ARG
46	BP	59	LEU
46	BP	61	ARG
46	BP	62	LEU
46	BP	71	VAL
46	BP	75	ILE
46	BP	85	LEU
46	BP	91	PHE
46	BP	101	VAL
46	BP	107	LYS
46	BP	110	TYR
46	BP	112	LEU
46	BP	115	LEU
46	BP	122	PRO
46	BP	123	LEU
46	BP	125	VAL
46	BP	130	PHE
46	BP	138	LEU
46	BP	144	GLU
46	BP	147	LEU
46	BP	148	LEU
47	BQ	9	TYR
47	BQ	17	LEU
47	BQ	22	LYS
47	BQ	26	TYR
47	BQ	29	PHE
47	BQ	43	THR
47	BQ	55	VAL
47	BQ	57	HIS
47	BQ	68	ILE
47	BQ	79	LEU
47	BQ	80	GLU
47	BQ	82	ARG
47	BQ	83	MET
47	BQ	87	LYS
47	BQ	91	GLU
47	BQ	106	VAL
47	BQ	119	ARG

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Mol	Chain	Res	Type
47	BQ	127	ILE
47	BQ	131	ILE
47	BQ	137	TYR
48	BR	2	ARG
48	BR	3	HIS
48	BR	5	LYS
48	BR	13	HIS
48	BR	18	LEU
48	BR	20	LEU
48	BR	29	LEU
48	BR	44	LEU
48	BR	49	ASP
48	BR	60	LEU
48	BR	65	LEU
48	BR	71	GLN
48	BR	74	LYS
48	BR	75	LEU
48	BR	79	LEU
48	BR	99	LYS
48	BR	103	ARG
48	BR	104	ARG
48	BR	113	LEU
48	BR	118	GLU
49	BS	11	LYS
49	BS	12	PHE
49	BS	13	ARG
49	BS	16	ASN
49	BS	36	TYR
49	BS	56	LEU
49	BS	61	ASN
49	BS	62	LYS
49	BS	71	ARG
49	BS	89	ARG
49	BS	92	TYR
49	BS	93	LYS
49	BS	95	HIS
49	BS	106	ARG
50	BT	1	MET
50	BT	3	ARG
50	BT	11	GLU
50	BT	12	SER
50	BT	13	ARG

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Mol	Chain	Res	Type
50	BT	14	TYR
50	BT	22	PHE
50	BT	28	VAL
50	BT	29	ARG
50	BT	38	ASN
50	BT	41	ARG
50	BT	44	ASP
50	BT	53	ARG
50	BT	58	ASN
50	BT	63	VAL
50	BT	65	LYS
50	BT	81	PRO
50	BT	87	ASP
50	BT	88	ILE
50	BT	93	ARG
50	BT	96	ARG
50	BT	99	LEU
50	BT	101	PHE
50	BT	102	ILE
50	BT	108	ARG
50	BT	110	ILE
50	BT	112	ARG
50	BT	115	ARG
50	BT	122	ASP
50	BT	123	GLN
50	BT	128	GLU
51	BU	3	ARG
51	BU	9	VAL
51	BU	14	HIS
51	BU	28	ARG
51	BU	30	LYS
51	BU	31	SER
51	BU	40	PHE
51	BU	49	HIS
51	BU	64	ARG
51	BU	69	CYS
51	BU	92	ARG
51	BU	93	LYS
52	BV	6	LYS
52	BV	13	ARG
52	BV	18	LEU
52	BV	19	LYS

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Mol	Chain	Res	Type
52	BV	21	ARG
52	BV	40	LEU
52	BV	66	ARG
52	BV	71	LEU
52	BV	78	LYS
52	BV	82	ARG
52	BV	83	ARG
52	BV	88	ARG
52	BV	94	LEU
52	BV	98	GLU
53	BW	11	ARG
53	BW	19	LEU
53	BW	20	VAL
53	BW	49	LYS
53	BW	51	LEU
53	BW	57	ASN
53	BW	61	ASN
53	BW	67	ASP
53	BW	69	LEU
53	BW	70	TYR
53	BW	75	TYR
53	BW	88	ARG
53	BW	107	LEU
54	BX	16	LYS
54	BX	26	TYR
54	BX	27	THR
54	BX	28	PHE
54	BX	29	TRP
54	BX	36	LYS
54	BX	40	LYS
54	BX	45	THR
54	BX	55	ASN
54	BX	59	VAL
54	BX	65	ARG
54	BX	70	LEU
54	BX	76	ARG
54	BX	77	LYS
54	BX	78	LYS
55	BY	2	ARG
55	BY	6	HIS
55	BY	8	LYS
55	BY	28	LYS

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Mol	Chain	Res	Type
55	BY	29	GLU
55	BY	31	LEU
55	BY	32	PRO
55	BY	38	ILE
55	BY	47	LYS
55	BY	56	PRO
55	BY	60	PHE
55	BY	62	GLU
55	BY	66	PRO
55	BY	71	LYS
55	BY	72	VAL
55	BY	75	ILE
55	BY	76	CYS
55	BY	89	PHE
55	BY	92	ASN
55	BY	96	ILE
55	BY	97	ARG
56	BZ	6	LYS
56	BZ	9	TYR
56	BZ	11	GLU
56	BZ	20	ARG
56	BZ	34	ASN
56	BZ	45	ASP
56	BZ	48	PHE
56	BZ	55	HIS
56	BZ	74	VAL
56	BZ	79	ARG
56	BZ	87	ASP
56	BZ	91	LEU
56	BZ	97	GLU
56	BZ	103	ARG
56	BZ	111	VAL
56	BZ	121	HIS
56	BZ	126	VAL
56	BZ	136	PHE
56	BZ	140	ASP
56	BZ	150	LEU
56	BZ	155	LEU
2	CB	17	PHE
2	CB	22	LYS
2	CB	31	TYR
2	CB	33	TYR

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Mol	Chain	Res	Type
2	CB	44	LEU
2	CB	56	ARG
2	CB	67	THR
2	CB	69	LEU
2	CB	70	PHE
2	CB	75	LYS
2	CB	80	ILE
2	CB	96	ARG
2	CB	97	TRP
2	CB	98	LEU
2	CB	102	LEU
2	CB	108	ILE
2	CB	115	LEU
2	CB	116	GLU
2	CB	137	ARG
2	CB	140	HIS
2	CB	158	LEU
2	CB	178	ARG
2	CB	187	LEU
2	CB	212	GLN
2	CB	222	ILE
3	CC	5	ILE
3	CC	12	LEU
3	CC	16	ARG
3	CC	27	LYS
3	CC	37	GLN
3	CC	43	LEU
3	CC	52	LEU
3	CC	101	LEU
3	CC	104	GLN
3	CC	120	VAL
3	CC	127	ARG
3	CC	131	ARG
3	CC	156	ARG
3	CC	164	ARG
3	CC	165	THR
3	CC	166	GLU
3	CC	167	TRP
3	CC	181	ASN
4	CD	3	ARG
4	CD	9	CYS
4	CD	21	LEU

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Mol	Chain	Res	Type
4	CD	26	CYS
4	CD	31	CYS
4	CD	33	MET
4	CD	42	GLN
4	CD	53	ASP
4	CD	59	ARG
4	CD	79	PHE
4	CD	96	LEU
4	CD	97	LEU
4	CD	110	PHE
4	CD	120	LEU
4	CD	122	ARG
4	CD	131	ARG
4	CD	132	ARG
4	CD	138	TYR
4	CD	158	ILE
4	CD	170	VAL
4	CD	181	MET
4	CD	194	LEU
4	CD	196	LEU
4	CD	200	GLU
5	CE	10	MET
5	CE	12	LEU
5	CE	16	THR
5	CE	20	GLN
5	CE	47	LYS
5	CE	60	TYR
5	CE	61	TYR
5	CE	68	GLU
5	CE	71	LEU
5	CE	73	ASN
5	CE	76	ILE
5	CE	90	VAL
5	CE	96	PRO
5	CE	101	ILE
5	CE	131	ILE
5	CE	149	GLU
6	CF	32	ASN
6	CF	75	LEU
6	CF	98	LEU
6	CF	100	ASN
7	CG	14	PRO

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Mol	Chain	Res	Type
7	CG	41	ARG
7	CG	54	THR
7	CG	57	GLU
7	CG	84	ASN
7	CG	88	PRO
7	CG	111	ARG
7	CG	112	PRO
7	CG	151	TYR
7	CG	156	TRP
8	CH	1	MET
8	CH	3	THR
8	CH	10	LEU
8	CH	18	ARG
8	CH	25	ASP
8	CH	35	ILE
8	CH	102	ARG
8	CH	119	LEU
8	CH	133	LEU
8	CH	136	GLU
8	CH	137	VAL
9	CI	4	TYR
9	CI	27	THR
9	CI	63	ILE
9	CI	66	ARG
9	CI	78	LYS
9	CI	95	LYS
9	CI	101	PHE
9	CI	104	ARG
9	CI	108	VAL
9	CI	113	LYS
9	CI	114	TYR
9	CI	121	ARG
9	CI	125	TYR
10	CJ	16	LEU
10	CJ	22	LYS
10	CJ	47	PHE
10	CJ	50	ILE
10	CJ	62	HIS
10	CJ	84	GLN
10	CJ	96	ILE
11	CK	18	ARG
11	CK	21	ILE

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Mol	Chain	Res	Type
11	CK	48	ILE
11	CK	91	ARG
11	CK	103	LEU
11	CK	117	ASN
11	CK	125	PHE
11	CK	126	ARG
11	CK	127	LYS
12	CL	19	ARG
12	CL	20	LYS
12	CL	21	LYS
12	CL	41	ARG
12	CL	48	PRO
12	CL	49	ASN
12	CL	53	ARG
12	CL	66	VAL
12	CL	84	LEU
12	CL	89	ARG
12	CL	102	ARG
13	CM	47	ASP
13	CM	48	LEU
13	CM	64	TRP
13	CM	82	MET
13	CM	93	ARG
13	CM	108	ARG
13	CM	120	LYS
14	CN	16	PHE
14	CN	26	ARG
14	CN	29	ARG
14	CN	61	TRP
15	CO	4	THR
15	CO	6	GLU
15	CO	10	LYS
15	CO	35	ARG
15	CO	39	LEU
15	CO	46	HIS
15	CO	47	LYS
15	CO	65	ARG
15	CO	82	ILE
16	CP	1	MET
16	CP	6	LEU
16	CP	26	ARG
16	CP	27	LYS

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Mol	Chain	Res	Type
16	CP	29	ASP
16	CP	38	TYR
16	CP	39	TYR
16	CP	48	TRP
16	CP	69	THR
16	CP	73	LEU
16	CP	82	GLN
17	CQ	36	ILE
17	CQ	55	ASP
17	CQ	56	VAL
17	CQ	78	GLU
17	CQ	81	ARG
17	CQ	89	LEU
18	CR	29	PHE
18	CR	31	LEU
18	CR	43	PHE
18	CR	46	GLU
18	CR	58	LEU
18	CR	74	ARG
19	CS	5	LEU
19	CS	6	LYS
19	CS	10	PHE
19	CS	13	ASP
19	CS	27	GLU
19	CS	34	TRP
19	CS	37	ARG
19	CS	43	GLU
19	CS	44	MET
19	CS	49	ILE
19	CS	53	ASN
19	CS	70	LYS
19	CS	78	ARG
20	CT	8	ARG
20	CT	13	LEU
20	CT	24	LEU
20	CT	26	ASN
20	CT	57	ARG
20	CT	62	LEU
20	CT	71	THR
20	CT	73	HIS
20	CT	74	LYS
20	CT	93	GLU

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Mol	Chain	Res	Type
20	CT	100	ILE
21	CU	24	ARG
25	CY	3	LEU
25	CY	10	THR
25	CY	18	LEU
25	CY	32	ARG
25	CY	85	ASP
25	CY	91	ASN
25	CY	92	PRO
25	CY	97	ASP
25	CY	113	ASP
25	CY	118	VAL
25	CY	132	ILE
25	CY	138	ASP
25	CY	168	PHE
25	CY	173	ASP
25	CY	183	ILE
26	D0	20	ARG
26	D0	38	VAL
26	D0	44	ARG
26	D0	64	ASP
27	D1	12	PRO
27	D1	16	ASN
27	D1	19	GLN
27	D1	20	ARG
27	D1	21	ARG
27	D1	35	THR
27	D1	39	LYS
27	D1	46	LEU
27	D1	47	GLN
27	D1	48	LYS
27	D1	49	VAL
27	D1	50	ARG
27	D1	52	ARG
27	D1	53	VAL
27	D1	56	GLN
27	D1	58	ILE
27	D1	69	LYS
27	D1	73	LEU
27	D1	83	GLU
27	D1	85	LEU
27	D1	89	GLU

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Mol	Chain	Res	Type
27	D1	90	ILE
27	D1	94	LEU
28	D2	12	GLU
28	D2	14	ARG
28	D2	22	GLU
28	D2	26	ARG
28	D2	30	ARG
28	D2	32	LEU
28	D2	33	MET
28	D2	37	PHE
28	D2	46	GLN
28	D2	53	LEU
28	D2	59	ARG
28	D2	61	LEU
29	D3	8	LEU
29	D3	26	LEU
29	D3	31	LEU
31	D5	4	HIS
31	D5	5	PRO
31	D5	22	HIS
31	D5	32	PRO
31	D5	43	HIS
31	D5	48	GLU
31	D5	56	LYS
31	D5	58	LEU
32	D6	10	LEU
32	D6	20	ASN
32	D6	31	PRO
32	D6	35	GLU
32	D6	37	ARG
32	D6	42	TRP
32	D6	43	CYS
32	D6	44	ARG
32	D6	51	GLU
33	D7	8	ASN
33	D7	24	THR
33	D7	36	GLN
33	D7	43	THR
33	D7	44	PRO
33	D7	46	VAL
33	D7	48	LYS
34	D8	31	HIS

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Mol	Chain	Res	Type
34	D8	40	GLU
34	D8	44	LYS
34	D8	46	ARG
34	D8	48	PHE
34	D8	52	LYS
34	D8	60	LEU
34	D8	64	TYR
37	DC	36	LYS
37	DC	49	ILE
37	DC	64	LEU
37	DC	66	HIS
38	DD	5	LYS
38	DD	10	THR
38	DD	14	ARG
38	DD	15	PHE
38	DD	24	ILE
38	DD	26	LYS
38	DD	31	LYS
38	DD	36	PRO
38	DD	46	GLN
38	DD	71	ASP
38	DD	91	ARG
38	DD	94	LEU
38	DD	95	LEU
38	DD	96	HIS
38	DD	131	LEU
38	DD	135	PHE
38	DD	165	ILE
38	DD	166	GLN
38	DD	173	VAL
38	DD	198	ASN
38	DD	211	ARG
38	DD	221	VAL
38	DD	228	PRO
38	DD	241	PRO
38	DD	255	LYS
38	DD	261	LYS
38	DD	271	ILE
39	DE	21	VAL
39	DE	33	VAL
39	DE	44	TYR
39	DE	61	ARG

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Mol	Chain	Res	Type
39	DE	64	LYS
39	DE	67	PHE
39	DE	69	LYS
39	DE	78	LEU
39	DE	79	ARG
39	DE	113	PHE
39	DE	118	LYS
39	DE	119	ARG
39	DE	122	PHE
39	DE	127	ASP
39	DE	134	ILE
39	DE	141	ILE
39	DE	143	ASN
39	DE	178	GLU
39	DE	181	LEU
39	DE	183	LEU
39	DE	184	VAL
39	DE	192	ASN
39	DE	202	LYS
39	DE	203	LYS
40	DF	6	VAL
40	DF	23	ASP
40	DF	24	LEU
40	DF	28	ILE
40	DF	46	ARG
40	DF	50	SER
40	DF	53	THR
40	DF	56	GLU
40	DF	65	TRP
40	DF	66	PRO
40	DF	67	GLN
40	DF	74	ARG
40	DF	77	ASP
40	DF	104	LYS
40	DF	112	MET
40	DF	116	ASP
40	DF	136	THR
40	DF	156	LEU
40	DF	164	ARG
40	DF	169	ASN
40	DF	197	ASP
40	DF	201	VAL

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Mol	Chain	Res	Type
40	DF	202	PHE
41	DG	16	ARG
41	DG	19	LEU
41	DG	21	ARG
41	DG	22	ARG
41	DG	23	PHE
41	DG	34	LEU
41	DG	43	LEU
41	DG	45	GLU
41	DG	51	ARG
41	DG	53	LEU
41	DG	58	GLN
41	DG	59	GLU
41	DG	76	SER
41	DG	77	ILE
41	DG	95	ARG
41	DG	121	ASN
41	DG	123	ASN
41	DG	125	PHE
41	DG	126	ASP
41	DG	131	TYR
41	DG	133	LEU
41	DG	136	ARG
41	DG	143	GLU
41	DG	150	ASP
41	DG	155	MET
41	DG	160	VAL
41	DG	170	ARG
41	DG	174	GLU
42	DH	41	MET
42	DH	53	GLU
42	DH	61	HIS
42	DH	68	THR
42	DH	83	TYR
42	DH	86	GLU
42	DH	89	ILE
42	DH	105	LEU
42	DH	123	PHE
42	DH	136	ILE
42	DH	143	GLN
42	DH	153	LYS
42	DH	157	TYR

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Mol	Chain	Res	Type
42	DH	158	HIS
42	DH	159	GLU
42	DH	170	ARG
43	DI	1	MET
43	DI	3	VAL
43	DI	5	LEU
43	DI	7	GLU
43	DI	10	GLU
43	DI	12	LEU
43	DI	25	TYR
43	DI	38	LEU
43	DI	51	ILE
43	DI	54	GLN
43	DI	64	GLU
43	DI	86	THR
43	DI	96	ASP
43	DI	118	LYS
43	DI	122	GLU
43	DI	123	LEU
43	DI	126	TYR
43	DI	130	TYR
43	DI	136	VAL
44	DN	4	TYR
44	DN	19	GLU
44	DN	34	LEU
44	DN	39	ARG
44	DN	42	TRP
44	DN	56	ASN
44	DN	62	VAL
44	DN	66	LYS
44	DN	82	LEU
44	DN	87	LEU
44	DN	94	HIS
44	DN	99	LEU
44	DN	109	LYS
44	DN	111	PRO
44	DN	123	TYR
44	DN	127	ASP
44	DN	128	HIS
44	DN	131	GLN
45	DO	1	MET
45	DO	7	TYR

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Mol	Chain	Res	Type
45	DO	9	GLU
45	DO	19	ILE
45	DO	24	VAL
45	DO	25	LEU
45	DO	32	TYR
45	DO	39	ILE
45	DO	45	GLU
45	DO	47	ILE
45	DO	48	PRO
45	DO	69	ILE
45	DO	73	ASP
45	DO	80	ASP
45	DO	82	ASN
45	DO	86	ILE
45	DO	87	ILE
45	DO	91	LEU
45	DO	102	VAL
45	DO	104	ARG
45	DO	115	VAL
46	DP	6	LEU
46	DP	13	ASN
46	DP	16	ARG
46	DP	18	ARG
46	DP	21	ARG
46	DP	32	THR
46	DP	35	HIS
46	DP	50	ARG
46	DP	59	LEU
46	DP	61	ARG
46	DP	62	LEU
46	DP	71	VAL
46	DP	75	ILE
46	DP	85	LEU
46	DP	91	PHE
46	DP	101	VAL
46	DP	107	LYS
46	DP	110	TYR
46	DP	112	LEU
46	DP	115	LEU
46	DP	122	PRO
46	DP	123	LEU
46	DP	125	VAL

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Mol	Chain	Res	Type
46	DP	130	PHE
46	DP	138	LEU
46	DP	144	GLU
46	DP	147	LEU
46	DP	148	LEU
47	DQ	9	TYR
47	DQ	17	LEU
47	DQ	22	LYS
47	DQ	26	TYR
47	DQ	29	PHE
47	DQ	43	THR
47	DQ	55	VAL
47	DQ	57	HIS
47	DQ	68	ILE
47	DQ	79	LEU
47	DQ	80	GLU
47	DQ	83	MET
47	DQ	87	LYS
47	DQ	91	GLU
47	DQ	106	VAL
47	DQ	127	ILE
47	DQ	131	ILE
47	DQ	137	TYR
48	DR	2	ARG
48	DR	3	HIS
48	DR	5	LYS
48	DR	13	HIS
48	DR	18	LEU
48	DR	20	LEU
48	DR	29	LEU
48	DR	44	LEU
48	DR	49	ASP
48	DR	60	LEU
48	DR	65	LEU
48	DR	71	GLN
48	DR	74	LYS
48	DR	75	LEU
48	DR	79	LEU
48	DR	99	LYS
48	DR	103	ARG
48	DR	104	ARG
48	DR	113	LEU

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Mol	Chain	Res	Type
48	DR	118	GLU
49	DS	11	LYS
49	DS	12	PHE
49	DS	13	ARG
49	DS	16	ASN
49	DS	36	TYR
49	DS	54	LEU
49	DS	56	LEU
49	DS	61	ASN
49	DS	62	LYS
49	DS	71	ARG
49	DS	89	ARG
49	DS	92	TYR
49	DS	93	LYS
49	DS	95	HIS
49	DS	106	ARG
50	DT	1	MET
50	DT	3	ARG
50	DT	11	GLU
50	DT	12	SER
50	DT	13	ARG
50	DT	14	TYR
50	DT	22	PHE
50	DT	28	VAL
50	DT	29	ARG
50	DT	38	ASN
50	DT	41	ARG
50	DT	44	ASP
50	DT	53	ARG
50	DT	58	ASN
50	DT	63	VAL
50	DT	65	LYS
50	DT	74	ARG
50	DT	81	PRO
50	DT	87	ASP
50	DT	93	ARG
50	DT	96	ARG
50	DT	99	LEU
50	DT	101	PHE
50	DT	102	ILE
50	DT	108	ARG
50	DT	110	ILE

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Mol	Chain	Res	Type
50	DT	112	ARG
50	DT	115	ARG
50	DT	122	ASP
50	DT	123	GLN
50	DT	128	GLU
51	DU	3	ARG
51	DU	9	VAL
51	DU	14	HIS
51	DU	28	ARG
51	DU	30	LYS
51	DU	31	SER
51	DU	49	HIS
51	DU	64	ARG
51	DU	69	CYS
51	DU	92	ARG
51	DU	93	LYS
52	DV	6	LYS
52	DV	13	ARG
52	DV	18	LEU
52	DV	19	LYS
52	DV	21	ARG
52	DV	40	LEU
52	DV	66	ARG
52	DV	71	LEU
52	DV	78	LYS
52	DV	82	ARG
52	DV	83	ARG
52	DV	88	ARG
52	DV	94	LEU
52	DV	98	GLU
53	DW	11	ARG
53	DW	19	LEU
53	DW	20	VAL
53	DW	49	LYS
53	DW	51	LEU
53	DW	57	ASN
53	DW	61	ASN
53	DW	67	ASP
53	DW	69	LEU
53	DW	70	TYR
53	DW	75	TYR
53	DW	88	ARG

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Mol	Chain	Res	Type
53	DW	107	LEU
54	DX	16	LYS
54	DX	26	TYR
54	DX	27	THR
54	DX	28	PHE
54	DX	29	TRP
54	DX	36	LYS
54	DX	40	LYS
54	DX	45	THR
54	DX	55	ASN
54	DX	59	VAL
54	DX	65	ARG
54	DX	66	LEU
54	DX	70	LEU
54	DX	76	ARG
54	DX	77	LYS
54	DX	78	LYS
55	DY	2	ARG
55	DY	6	HIS
55	DY	8	LYS
55	DY	28	LYS
55	DY	29	GLU
55	DY	31	LEU
55	DY	32	PRO
55	DY	38	ILE
55	DY	47	LYS
55	DY	56	PRO
55	DY	60	PHE
55	DY	62	GLU
55	DY	66	PRO
55	DY	71	LYS
55	DY	72	VAL
55	DY	75	ILE
55	DY	76	CYS
55	DY	89	PHE
55	DY	92	ASN
55	DY	96	ILE
55	DY	97	ARG
56	DZ	5	LEU
56	DZ	6	LYS
56	DZ	9	TYR
56	DZ	10	ARG

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Mol	Chain	Res	Type
56	DZ	11	GLU
56	DZ	19	ARG
56	DZ	24	LEU
56	DZ	39	VAL
56	DZ	40	ASP
56	DZ	41	LEU
56	DZ	42	VAL
56	DZ	45	ASP
56	DZ	59	LEU
56	DZ	71	VAL
56	DZ	72	ARG
56	DZ	74	VAL
56	DZ	76	LEU
56	DZ	79	ARG
56	DZ	84	GLU
56	DZ	91	LEU
56	DZ	103	ARG
56	DZ	108	PRO
56	DZ	119	GLU
56	DZ	121	HIS
56	DZ	123	ASP
56	DZ	136	PHE
56	DZ	144	LEU
56	DZ	148	ASP
56	DZ	150	LEU
56	DZ	163	LEU
56	DZ	166	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (303) such sidechains are listed below:

Mol	Chain	Res	Type
2	AB	37	ASN
2	AB	78	GLN
2	AB	135	GLN
2	AB	146	GLN
2	AB	204	ASN
2	AB	212	GLN
3	AC	28	GLN
3	AC	31	HIS
3	AC	69	HIS
3	AC	98	ASN
3	AC	123	GLN

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Mol	Chain	Res	Type
3	AC	170	GLN
3	AC	181	ASN
4	AD	45	GLN
4	AD	62	GLN
4	AD	119	GLN
4	AD	161	ASN
4	AD	201	GLN
5	AE	56	GLN
5	AE	73	ASN
5	AE	78	HIS
5	AE	127	ASN
6	AF	7	ASN
6	AF	18	GLN
6	AF	27	GLN
6	AF	32	ASN
6	AF	94	GLN
6	AF	100	ASN
7	AG	11	GLN
7	AG	13	GLN
7	AG	28	ASN
7	AG	68	ASN
7	AG	84	ASN
7	AG	106	GLN
7	AG	109	ASN
7	AG	148	ASN
8	AH	15	ASN
8	AH	78	GLN
8	AH	82	HIS
9	AI	23	ASN
9	AI	124	GLN
10	AJ	13	HIS
11	AK	26	ASN
11	AK	104	GLN
11	AK	117	ASN
12	AL	8	ASN
12	AL	9	GLN
12	AL	49	ASN
12	AL	75	HIS
12	AL	99	HIS
13	AM	40	ASN
13	AM	77	ASN
15	AO	13	GLN

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Mol	Chain	Res	Type
15	AO	37	ASN
15	AO	46	HIS
15	AO	62	GLN
16	AP	13	HIS
16	AP	76	GLN
16	AP	82	GLN
17	AQ	16	GLN
18	AR	36	ASN
19	AS	47	HIS
19	AS	53	ASN
20	AT	16	HIS
20	AT	18	GLN
20	AT	26	ASN
25	AY	34	ASN
25	AY	53	ASN
25	AY	91	ASN
25	AY	174	GLN
26	B0	29	GLN
26	B0	35	ASN
27	B1	19	GLN
27	B1	42	GLN
27	B1	45	ASN
27	B1	66	HIS
28	B2	46	GLN
28	B2	56	GLN
29	B3	19	GLN
29	B3	46	ASN
31	B5	23	HIS
31	B5	43	HIS
32	B6	20	ASN
32	B6	26	ASN
32	B6	32	ASN
33	B7	6	GLN
33	B7	8	ASN
34	B8	35	GLN
37	BC	44	HIS
37	BC	56	GLN
37	BC	57	ASN
38	BD	58	HIS
38	BD	126	GLN
38	BD	166	GLN
38	BD	186	HIS

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Mol	Chain	Res	Type
38	BD	198	ASN
39	BE	35	GLN
39	BE	48	GLN
39	BE	54	GLN
39	BE	55	ASN
39	BE	132	HIS
39	BE	192	ASN
40	BF	8	GLN
40	BF	69	HIS
40	BF	75	HIS
40	BF	160	ASN
40	BF	169	ASN
41	BG	27	ASN
41	BG	41	GLN
41	BG	123	ASN
42	BH	74	ASN
43	BI	11	ASN
43	BI	43	ASN
43	BI	105	HIS
44	BN	56	ASN
44	BN	101	HIS
44	BN	128	HIS
45	BO	3	GLN
45	BO	13	ASN
46	BP	13	ASN
46	BP	35	HIS
47	BQ	12	GLN
47	BQ	13	GLN
47	BQ	123	HIS
47	BQ	141	GLN
48	BR	11	ASN
48	BR	13	HIS
48	BR	61	HIS
48	BR	91	GLN
49	BS	16	ASN
49	BS	61	ASN
49	BS	84	GLN
50	BT	38	ASN
50	BT	55	ASN
50	BT	58	ASN
50	BT	84	GLN
50	BT	90	GLN

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Mol	Chain	Res	Type
50	BT	123	GLN
51	BU	49	HIS
51	BU	66	ASN
51	BU	81	HIS
51	BU	94	ASN
51	BU	117	GLN
52	BV	11	GLN
52	BV	89	GLN
53	BW	40	ASN
53	BW	57	ASN
54	BX	31	HIS
54	BX	41	ASN
54	BX	55	ASN
56	BZ	73	GLN
56	BZ	118	GLN
56	BZ	132	ASN
2	CB	37	ASN
2	CB	78	GLN
2	CB	135	GLN
2	CB	146	GLN
2	CB	204	ASN
2	CB	212	GLN
3	CC	28	GLN
3	CC	31	HIS
3	CC	69	HIS
3	CC	98	ASN
3	CC	123	GLN
3	CC	170	GLN
3	CC	181	ASN
4	CD	62	GLN
4	CD	119	GLN
4	CD	161	ASN
4	CD	201	GLN
5	CE	56	GLN
5	CE	73	ASN
5	CE	78	HIS
5	CE	127	ASN
6	CF	7	ASN
6	CF	27	GLN
6	CF	32	ASN
6	CF	94	GLN
6	CF	100	ASN

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Mol	Chain	Res	Type
7	CG	11	GLN
7	CG	13	GLN
7	CG	28	ASN
7	CG	68	ASN
7	CG	84	ASN
7	CG	106	GLN
7	CG	109	ASN
7	CG	148	ASN
8	CH	15	ASN
8	CH	78	GLN
8	CH	82	HIS
9	CI	23	ASN
9	CI	124	GLN
10	CJ	13	HIS
11	CK	104	GLN
11	CK	117	ASN
12	CL	8	ASN
12	CL	9	GLN
12	CL	49	ASN
12	CL	75	HIS
12	CL	99	HIS
13	CM	40	ASN
13	CM	77	ASN
13	CM	101	GLN
15	CO	13	GLN
15	CO	37	ASN
15	CO	46	HIS
16	CP	13	HIS
16	CP	76	GLN
16	CP	82	GLN
17	CQ	16	GLN
18	CR	36	ASN
19	CS	47	HIS
19	CS	53	ASN
20	CT	16	HIS
20	CT	18	GLN
20	CT	26	ASN
25	CY	34	ASN
25	CY	91	ASN
25	CY	174	GLN
26	D0	29	GLN
26	D0	35	ASN

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Mol	Chain	Res	Type
27	D1	16	ASN
27	D1	19	GLN
27	D1	56	GLN
28	D2	56	GLN
29	D3	19	GLN
29	D3	46	ASN
31	D5	23	HIS
31	D5	43	HIS
32	D6	20	ASN
32	D6	26	ASN
32	D6	32	ASN
33	D7	6	GLN
33	D7	8	ASN
34	D8	35	GLN
37	DC	44	HIS
37	DC	56	GLN
37	DC	57	ASN
38	DD	58	HIS
38	DD	126	GLN
38	DD	166	GLN
38	DD	186	HIS
38	DD	198	ASN
38	DD	253	GLN
39	DE	35	GLN
39	DE	48	GLN
39	DE	54	GLN
39	DE	55	ASN
39	DE	132	HIS
39	DE	192	ASN
40	DF	8	GLN
40	DF	69	HIS
40	DF	75	HIS
40	DF	160	ASN
40	DF	169	ASN
41	DG	27	ASN
41	DG	40	ASN
41	DG	66	GLN
41	DG	121	ASN
41	DG	130	ASN
41	DG	132	ASN
42	DH	74	ASN
42	DH	147	ASN

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Mol	Chain	Res	Type
43	DI	11	ASN
43	DI	43	ASN
43	DI	105	HIS
44	DN	56	ASN
44	DN	101	HIS
44	DN	128	HIS
45	DO	3	GLN
45	DO	13	ASN
46	DP	13	ASN
47	DQ	12	GLN
47	DQ	13	GLN
47	DQ	123	HIS
47	DQ	141	GLN
48	DR	11	ASN
48	DR	13	HIS
48	DR	61	HIS
48	DR	91	GLN
49	DS	16	ASN
49	DS	61	ASN
49	DS	84	GLN
50	DT	38	ASN
50	DT	55	ASN
50	DT	58	ASN
50	DT	84	GLN
50	DT	90	GLN
50	DT	123	GLN
51	DU	49	HIS
51	DU	66	ASN
51	DU	81	HIS
51	DU	94	ASN
51	DU	117	GLN
52	DV	11	GLN
52	DV	89	GLN
53	DW	40	ASN
53	DW	57	ASN
54	DX	31	HIS
54	DX	41	ASN
54	DX	55	ASN
55	DY	92	ASN
56	DZ	55	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1503/1522 (98%)	210 (13%)	31 (2%)
1	CA	1503/1522 (98%)	214 (14%)	29 (1%)
22	AV	16/76 (21%)	0	0
22	CV	16/76 (21%)	0	0
23	AW	76/77 (98%)	26 (34%)	0
23	CW	76/77 (98%)	20 (26%)	0
24	AX	10/31 (32%)	3 (30%)	0
24	CX	10/31 (32%)	2 (20%)	0
35	BA	2766/2782 (99%)	577 (20%)	57 (2%)
35	DA	2766/2782 (99%)	580 (20%)	57 (2%)
36	BB	118/122 (96%)	19 (16%)	1 (0%)
36	DB	118/122 (96%)	19 (16%)	1 (0%)
All	All	8978/9220 (97%)	1670 (18%)	176 (1%)

All (1670) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	9	G
1	AA	31	G
1	AA	32	A
1	AA	39	G
1	AA	47	C
1	AA	48	C
1	AA	51	A
1	AA	61	G
1	AA	65	U
1	AA	80	G
1	AA	81	U
1	AA	90	U
1	AA	97	G
1	AA	98	G
1	AA	116	A
1	AA	120	A
1	AA	121	C
1	AA	131	C
1	AA	146	G
1	AA	150	C
1	AA	172	A
1	AA	189(F)	U
1	AA	189(H)	G
1	AA	195	A
1	AA	197	A
1	AA	204	U

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Mol	Chain	Res	Type
1	AA	244	U
1	AA	247	G
1	AA	251	G
1	AA	266	G
1	AA	267	C
1	AA	274	A
1	AA	279	A
1	AA	289	G
1	AA	316	G
1	AA	328	C
1	AA	329	A
1	AA	332	G
1	AA	345	C
1	AA	352	C
1	AA	353	A
1	AA	354	G
1	AA	367	U
1	AA	372	C
1	AA	397	A
1	AA	412	A
1	AA	413	G
1	AA	414	A
1	AA	422	C
1	AA	423	G
1	AA	428	G
1	AA	429	U
1	AA	430	A
1	AA	435	C
1	AA	436	C
1	AA	437	U
1	AA	442	C
1	AA	452	A
1	AA	461	A
1	AA	482	A
1	AA	484	G
1	AA	485	G
1	AA	496	A
1	AA	498	U
1	AA	509	A
1	AA	510	A
1	AA	511	C
1	AA	518	C

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Mol	Chain	Res	Type
1	AA	524	G
1	AA	527	G
1	AA	531	U
1	AA	532	A
1	AA	533	A
1	AA	534	U
1	AA	536	C
1	AA	547	A
1	AA	559	A
1	AA	561	U
1	AA	562	C
1	AA	564	C
1	AA	566	G
1	AA	570	G
1	AA	572	A
1	AA	573	A
1	AA	575	G
1	AA	576	G
1	AA	577	G
1	AA	596	C
1	AA	631	G
1	AA	632	A
1	AA	653	A
1	AA	687	A
1	AA	688	G
1	AA	731	G
1	AA	749	C
1	AA	755	G
1	AA	793	U
1	AA	794	A
1	AA	803	G
1	AA	816	A
1	AA	817	C
1	AA	819	A
1	AA	820	U
1	AA	821	G
1	AA	828	A
1	AA	833	U
1	AA	839	U
1	AA	840	C
1	AA	841	U
1	AA	848	C

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Mol	Chain	Res	Type
1	AA	859	A
1	AA	874	G
1	AA	902	G
1	AA	914	A
1	AA	926	G
1	AA	927	G
1	AA	934	C
1	AA	935	A
1	AA	960	U
1	AA	966	G
1	AA	968	A
1	AA	969	A
1	AA	974	A
1	AA	975	A
1	AA	976	G
1	AA	977	A
1	AA	978	A
1	AA	980	C
1	AA	991	U
1	AA	992	U
1	AA	993	G
1	AA	1026	G
1	AA	1050	G
1	AA	1053	G
1	AA	1054	C
1	AA	1064	G
1	AA	1065	U
1	AA	1066	C
1	AA	1068	G
1	AA	1085	U
1	AA	1094	G
1	AA	1095	U
1	AA	1101	A
1	AA	1117	G
1	AA	1118	C
1	AA	1123	A
1	AA	1124	G
1	AA	1125	U
1	AA	1126	U
1	AA	1129	C
1	AA	1130	A
1	AA	1131	G

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Mol	Chain	Res	Type
1	AA	1137	C
1	AA	1138	G
1	AA	1139	G
1	AA	1146	A
1	AA	1147	C
1	AA	1152	A
1	AA	1159	U
1	AA	1184	G
1	AA	1187	G
1	AA	1190	G
1	AA	1196	U
1	AA	1197	G
1	AA	1212	U
1	AA	1213	A
1	AA	1225	A
1	AA	1227	A
1	AA	1238	A
1	AA	1249	C
1	AA	1256	A
1	AA	1257	U
1	AA	1278	U
1	AA	1280	A
1	AA	1281	U
1	AA	1282	C
1	AA	1286	A
1	AA	1287	A
1	AA	1294	G
1	AA	1297	C
1	AA	1300	G
1	AA	1301	U
1	AA	1302	U
1	AA	1305	G
1	AA	1317	C
1	AA	1320	C
1	AA	1322	C
1	AA	1325	C
1	AA	1331	G
1	AA	1335	C
1	AA	1346	A
1	AA	1347	G
1	AA	1364	U
1	AA	1398	A

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Mol	Chain	Res	Type
1	AA	1419	G
1	AA	1432	G
1	AA	1442(A)	G
1	AA	1442(B)	A
1	AA	1443	G
1	AA	1452	C
1	AA	1456	G
1	AA	1487	G
1	AA	1492	A
1	AA	1499	A
1	AA	1504	G
1	AA	1505	G
1	AA	1506	U
1	AA	1517	G
1	AA	1529	G
1	AA	1530	G
23	AW	3	C
23	AW	4	G
23	AW	5	G
23	AW	8	U
23	AW	9	G
23	AW	17	C
23	AW	18	U
23	AW	19	G
23	AW	20	G
23	AW	21	U
23	AW	22	A
23	AW	24	C
23	AW	31	G
23	AW	33	C
23	AW	38	A
23	AW	39	A
23	AW	41	C
23	AW	48	U
23	AW	49	C
23	AW	50	G
23	AW	54	G
23	AW	61	U
23	AW	62	C
23	AW	74	A
23	AW	75	C
23	AW	77	A

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Mol	Chain	Res	Type
24	AX	13	A
24	AX	17	U
24	AX	22	U
35	BA	10	G
35	BA	17	G
35	BA	28	A
35	BA	34	C
35	BA	35	G
35	BA	45	C
35	BA	49	A
35	BA	50	U
35	BA	51	G
35	BA	55	G
35	BA	68	G
35	BA	69	C
35	BA	71	A
35	BA	72	U
35	BA	73	A
35	BA	75	G
35	BA	83	G
35	BA	84	A
35	BA	85	G
35	BA	88	G
35	BA	90	U
35	BA	94	C
35	BA	95	G
35	BA	99	U
35	BA	100	G
35	BA	102	G
35	BA	118	A
35	BA	120	U
35	BA	132	G
35	BA	139	G
35	BA	139(A)	G
35	BA	140	G
35	BA	142	A
35	BA	142(A)	C
35	BA	143(A)	C
35	BA	146	G
35	BA	149	A
35	BA	157	U
35	BA	158	U

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Mol	Chain	Res	Type
35	BA	174	C
35	BA	182	A
35	BA	196	A
35	BA	197	A
35	BA	199	A
35	BA	205	G
35	BA	215	G
35	BA	216	A
35	BA	221	A
35	BA	222	A
35	BA	228	A
35	BA	229	A
35	BA	248	G
35	BA	249	C
35	BA	252	G
35	BA	261	G
35	BA	271(K)	U
35	BA	271(L)	U
35	BA	271(M)	G
35	BA	271(N)	U
35	BA	271(O)	C
35	BA	271(P)	C
35	BA	271(T)	C
35	BA	271(Y)	U
35	BA	272	G
35	BA	272(B)	G
35	BA	272(H)	C
35	BA	274	G
35	BA	275	G
35	BA	283	A
35	BA	284	U
35	BA	286	C
35	BA	287	C
35	BA	292	C
35	BA	310	A
35	BA	311	A
35	BA	329	G
35	BA	330	A
35	BA	332	A
35	BA	333	G
35	BA	335	C
35	BA	343	C

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Mol	Chain	Res	Type
35	BA	349	G
35	BA	352	G
35	BA	353	G
35	BA	356	G
35	BA	362	U
35	BA	363(F)	A
35	BA	365	C
35	BA	372	G
35	BA	386	G
35	BA	387	U
35	BA	388	G
35	BA	405	U
35	BA	406	G
35	BA	411	G
35	BA	416	C
35	BA	418	G
35	BA	428	A
35	BA	444	C
35	BA	448	U
35	BA	451	C
35	BA	456	C
35	BA	457	A
35	BA	470	A
35	BA	475	U
35	BA	481	G
35	BA	482	A
35	BA	494	G
35	BA	505	A
35	BA	508	G
35	BA	509	C
35	BA	512	G
35	BA	525	U
35	BA	529	A
35	BA	531	C
35	BA	532	A
35	BA	533	G
35	BA	537	C
35	BA	542	C
35	BA	543	C
35	BA	547	A
35	BA	548	A
35	BA	549	G

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Mol	Chain	Res	Type
35	BA	551	G
35	BA	556	G
35	BA	562	U
35	BA	563	G
35	BA	572	A
35	BA	573	G
35	BA	575	A
35	BA	586	A
35	BA	588	U
35	BA	604	G
35	BA	607	U
35	BA	614(A)	U
35	BA	614(B)	G
35	BA	614(C)	A
35	BA	615	G
35	BA	620	G
35	BA	622	G
35	BA	627	A
35	BA	637	A
35	BA	645	C
35	BA	646	A
35	BA	650	C
35	BA	652	C
35	BA	656	G
35	BA	670	A
35	BA	671	C
35	BA	686	G
35	BA	687	C
35	BA	708	C
35	BA	717	G
35	BA	722	A
35	BA	730	C
35	BA	753	C
35	BA	764	A
35	BA	765	G
35	BA	776	G
35	BA	782	A
35	BA	784	A
35	BA	785	G
35	BA	790	C
35	BA	791	C
35	BA	792	G

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Mol	Chain	Res	Type
35	BA	805	G
35	BA	812	C
35	BA	819	A
35	BA	827	U
35	BA	828	U
35	BA	846	C
35	BA	848	G
35	BA	857	C
35	BA	859	G
35	BA	864	G
35	BA	866	A
35	BA	878	A
35	BA	892	G
35	BA	896	A
35	BA	897	C
35	BA	898	C
35	BA	899	A
35	BA	904	C
35	BA	910	A
35	BA	917	A
35	BA	926	A
35	BA	932	G
35	BA	941	A
35	BA	945	A
35	BA	946	G
35	BA	955	C
35	BA	956	G
35	BA	957	A
35	BA	958	U
35	BA	959	A
35	BA	961	C
35	BA	964	C
35	BA	965	C
35	BA	973	A
35	BA	974	G
35	BA	975	C
35	BA	980	A
35	BA	983	A
35	BA	985	C
35	BA	990	A
35	BA	991	C
35	BA	996	A

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Mol	Chain	Res	Type
35	BA	1011	G
35	BA	1012	U
35	BA	1013	C
35	BA	1017	G
35	BA	1020	A
35	BA	1022	G
35	BA	1023	U
35	BA	1025	G
35	BA	1026	U
35	BA	1033	U
35	BA	1039	G
35	BA	1040	C
35	BA	1041	C
35	BA	1042	G
35	BA	1044	G
35	BA	1045	A
35	BA	1047	G
35	BA	1049	C
35	BA	1052	C
35	BA	1106	A
35	BA	1110	G
35	BA	1111	A
35	BA	1112	G
35	BA	1113	U
35	BA	1115	G
35	BA	1129	A
35	BA	1130	U
35	BA	1135	C
35	BA	1136	G
35	BA	1143	A
35	BA	1155	A
35	BA	1156	A
35	BA	1159	U
35	BA	1173	G
35	BA	1174	A
35	BA	1175	U
35	BA	1176	G
35	BA	1177	A
35	BA	1178	C
35	BA	1194	A
35	BA	1195	G
35	BA	1205	U

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Mol	Chain	Res	Type
35	BA	1210	A
35	BA	1211	U
35	BA	1212	G
35	BA	1218	C
35	BA	1221	C
35	BA	1241	A
35	BA	1247	A
35	BA	1248	G
35	BA	1249	U
35	BA	1251	C
35	BA	1253	A
35	BA	1254	A
35	BA	1255	U
35	BA	1256	G
35	BA	1265	A
35	BA	1271	G
35	BA	1272	A
35	BA	1276	A
35	BA	1281	G
35	BA	1300	U
35	BA	1301	A
35	BA	1302	A
35	BA	1314	C
35	BA	1319	G
35	BA	1329	U
35	BA	1330	C
35	BA	1332	G
35	BA	1345	C
35	BA	1349	A
35	BA	1359	A
35	BA	1365	A
35	BA	1368	G
35	BA	1379	A
35	BA	1380	G
35	BA	1385	G
35	BA	1386	C
35	BA	1390	U
35	BA	1407	C
35	BA	1412	A
35	BA	1416	G
35	BA	1417	C
35	BA	1420	U

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Mol	Chain	Res	Type
35	BA	1421	G
35	BA	1428	C
35	BA	1437	C
35	BA	1445	A
35	BA	1449	A
35	BA	1450	G
35	BA	1460	A
35	BA	1461	G
35	BA	1467	C
35	BA	1471	A
35	BA	1475	G
35	BA	1478	G
35	BA	1482	G
35	BA	1484	G
35	BA	1485	G
35	BA	1490	A
35	BA	1493	C
35	BA	1494	A
35	BA	1495	A
35	BA	1497	U
35	BA	1498	C
35	BA	1502	C
35	BA	1505	C
35	BA	1508	A
35	BA	1509	C
35	BA	1509(A)	A
35	BA	1528(A)	A
35	BA	1529	G
35	BA	1530	C
35	BA	1532	C
35	BA	1533	G
35	BA	1543	C
35	BA	1545	A
35	BA	1547	C
35	BA	1554	A
35	BA	1558	A
35	BA	1559	G
35	BA	1569	A
35	BA	1578	U
35	BA	1579	A
35	BA	1584	C
35	BA	1586	A

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Mol	Chain	Res	Type
35	BA	1587	A
35	BA	1591	G
35	BA	1594	G
35	BA	1595	G
35	BA	1603	A
35	BA	1608	A
35	BA	1609	A
35	BA	1610	A
35	BA	1613	G
35	BA	1614	A
35	BA	1615	C
35	BA	1616	A
35	BA	1617	C
35	BA	1635	G
35	BA	1640	C
35	BA	1648	C
35	BA	1654	A
35	BA	1674	G
35	BA	1678	G
35	BA	1694	C
35	BA	1696	G
35	BA	1707	G
35	BA	1718	G
35	BA	1739	U
35	BA	1741	A
35	BA	1742	G
35	BA	1746	G
35	BA	1748	G
35	BA	1763	G
35	BA	1764	G
35	BA	1773	A
35	BA	1780	A
35	BA	1791	A
35	BA	1799	G
35	BA	1800	C
35	BA	1816	G
35	BA	1820	U
35	BA	1835	G
35	BA	1847	A
35	BA	1854	A
35	BA	1858	G
35	BA	1865	G

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Mol	Chain	Res	Type
35	BA	1866	C
35	BA	1877	A
35	BA	1878	G
35	BA	1880	C
35	BA	1882	C
35	BA	1885	A
35	BA	1888	G
35	BA	1900	A
35	BA	1903	G
35	BA	1906	G
35	BA	1912	A
35	BA	1913	A
35	BA	1929	G
35	BA	1936	A
35	BA	1938	A
35	BA	1955	U
35	BA	1963	U
35	BA	1964	G
35	BA	1965	C
35	BA	1967	C
35	BA	1969	A
35	BA	1971	A
35	BA	1972	A
35	BA	1987	G
35	BA	1988	C
35	BA	1991	U
35	BA	1993	U
35	BA	2023	G
35	BA	2031	A
35	BA	2033	A
35	BA	2034	U
35	BA	2036	C
35	BA	2043	C
35	BA	2051	A
35	BA	2055	C
35	BA	2056	G
35	BA	2060	A
35	BA	2061	G
35	BA	2062	A
35	BA	2069	G
35	BA	2093	G
35	BA	2099	U

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Mol	Chain	Res	Type
35	BA	2103	C
35	BA	2104	G
35	BA	2108	C
35	BA	2110	G
35	BA	2111	C
35	BA	2112	G
35	BA	2116	G
35	BA	2117	A
35	BA	2118	U
35	BA	2119	A
35	BA	2120	G
35	BA	2122	U
35	BA	2127	G
35	BA	2128	C
35	BA	2163	C
35	BA	2164	C
35	BA	2165	G
35	BA	2166	G
35	BA	2169	A
35	BA	2171	A
35	BA	2172	U
35	BA	2173	A
35	BA	2176	A
35	BA	2179	C
35	BA	2185	C
35	BA	2187	G
35	BA	2190	G
35	BA	2198	A
35	BA	2199	A
35	BA	2200	C
35	BA	2207	G
35	BA	2208	A
35	BA	2218	U
35	BA	2225	A
35	BA	2226	C
35	BA	2238	G
35	BA	2239	G
35	BA	2243	U
35	BA	2251	G
35	BA	2263	C
35	BA	2275	C
35	BA	2283	C

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Mol	Chain	Res	Type
35	BA	2285	C
35	BA	2287	A
35	BA	2290	G
35	BA	2294	C
35	BA	2297	C
35	BA	2305	A
35	BA	2307	G
35	BA	2308	G
35	BA	2311	A
35	BA	2313	C
35	BA	2316	C
35	BA	2319	G
35	BA	2320	A
35	BA	2325	G
35	BA	2334	G
35	BA	2336	A
35	BA	2345	G
35	BA	2347	C
35	BA	2349	G
35	BA	2358	G
35	BA	2383	G
35	BA	2385	C
35	BA	2387	U
35	BA	2388	A
35	BA	2398	U
35	BA	2399	G
35	BA	2400	G
35	BA	2402	C
35	BA	2408	U
35	BA	2423	U
35	BA	2425	A
35	BA	2427	C
35	BA	2428	G
35	BA	2429	G
35	BA	2430	A
35	BA	2431	U
35	BA	2439	A
35	BA	2441	C
35	BA	2447	G
35	BA	2448	A
35	BA	2465	C
35	BA	2469	A

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Mol	Chain	Res	Type
35	BA	2470	G
35	BA	2474	C
35	BA	2476	A
35	BA	2478	A
35	BA	2482	G
35	BA	2483	C
35	BA	2490	G
35	BA	2491	U
35	BA	2498	C
35	BA	2502	G
35	BA	2505	G
35	BA	2518	A
35	BA	2520	C
35	BA	2523	G
35	BA	2529	G
35	BA	2534	A
35	BA	2543	G
35	BA	2554	U
35	BA	2559	C
35	BA	2566	A
35	BA	2567	G
35	BA	2571	C
35	BA	2572	A
35	BA	2573	C
35	BA	2582	G
35	BA	2585	U
35	BA	2586	C
35	BA	2602	A
35	BA	2609	U
35	BA	2610	C
35	BA	2611	U
35	BA	2612	C
35	BA	2615	U
35	BA	2630	G
35	BA	2640	G
35	BA	2646	C
35	BA	2655	G
35	BA	2673	G
35	BA	2675	A
35	BA	2682	U
35	BA	2690	C
35	BA	2691	C

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Mol	Chain	Res	Type
35	BA	2702	U
35	BA	2703	C
35	BA	2712(A)	A
35	BA	2713	A
35	BA	2720	U
35	BA	2726	U
35	BA	2733	A
35	BA	2752	C
35	BA	2754	U
35	BA	2759	G
35	BA	2762	G
35	BA	2763	G
35	BA	2765	A
35	BA	2778	A
35	BA	2779	U
35	BA	2781	A
35	BA	2789	C
35	BA	2790	A
35	BA	2791	C
35	BA	2796	U
35	BA	2801(A)	A
35	BA	2802	G
35	BA	2803	C
35	BA	2808	U
35	BA	2821	A
35	BA	2823	A
35	BA	2833	G
35	BA	2834	G
35	BA	2835	A
35	BA	2849	U
35	BA	2860	A
35	BA	2864	G
35	BA	2872	G
35	BA	2893	G
36	BB	8	U
36	BB	13	A
36	BB	15	A
36	BB	16	G
36	BB	22	U
36	BB	24	G
36	BB	25	A
36	BB	27	C

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Mol	Chain	Res	Type
36	BB	33	G
36	BB	41	U
36	BB	42	C
36	BB	45	A
36	BB	47	C
36	BB	53	A
36	BB	67	G
36	BB	73	A
36	BB	75	G
36	BB	88	C
36	BB	110	G
1	CA	9	G
1	CA	31	G
1	CA	32	A
1	CA	39	G
1	CA	47	C
1	CA	48	C
1	CA	51	A
1	CA	61	G
1	CA	65	U
1	CA	80	G
1	CA	81	U
1	CA	90	U
1	CA	97	G
1	CA	98	G
1	CA	116	A
1	CA	120	A
1	CA	121	C
1	CA	129(A)	G
1	CA	131	C
1	CA	146	G
1	CA	150	C
1	CA	172	A
1	CA	189(F)	U
1	CA	189(H)	G
1	CA	195	A
1	CA	197	A
1	CA	204	U
1	CA	244	U
1	CA	247	G
1	CA	251	G
1	CA	266	G

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Mol	Chain	Res	Type
1	CA	267	C
1	CA	274	A
1	CA	279	A
1	CA	289	G
1	CA	316	G
1	CA	328	C
1	CA	329	A
1	CA	332	G
1	CA	345	C
1	CA	352	C
1	CA	353	A
1	CA	354	G
1	CA	367	U
1	CA	372	C
1	CA	397	A
1	CA	412	A
1	CA	413	G
1	CA	414	A
1	CA	422	C
1	CA	423	G
1	CA	428	G
1	CA	429	U
1	CA	430	A
1	CA	435	C
1	CA	436	C
1	CA	437	U
1	CA	442	C
1	CA	452	A
1	CA	461	A
1	CA	482	A
1	CA	484	G
1	CA	485	G
1	CA	496	A
1	CA	498	U
1	CA	509	A
1	CA	510	A
1	CA	511	C
1	CA	518	C
1	CA	524	G
1	CA	527	G
1	CA	531	U
1	CA	532	A

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Mol	Chain	Res	Type
1	CA	533	A
1	CA	534	U
1	CA	536	C
1	CA	547	A
1	CA	559	A
1	CA	561	U
1	CA	562	C
1	CA	564	C
1	CA	566	G
1	CA	572	A
1	CA	573	A
1	CA	575	G
1	CA	576	G
1	CA	577	G
1	CA	596	C
1	CA	631	G
1	CA	632	A
1	CA	653	A
1	CA	687	A
1	CA	688	G
1	CA	731	G
1	CA	748	C
1	CA	749	C
1	CA	755	G
1	CA	793	U
1	CA	794	A
1	CA	803	G
1	CA	816	A
1	CA	817	C
1	CA	819	A
1	CA	820	U
1	CA	821	G
1	CA	828	A
1	CA	833	U
1	CA	839	U
1	CA	840	C
1	CA	841	U
1	CA	848	C
1	CA	859	A
1	CA	874	G
1	CA	902	G
1	CA	914	A

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Mol	Chain	Res	Type
1	CA	926	G
1	CA	927	G
1	CA	934	C
1	CA	935	A
1	CA	960	U
1	CA	966	G
1	CA	968	A
1	CA	969	A
1	CA	974	A
1	CA	975	A
1	CA	976	G
1	CA	977	A
1	CA	978	A
1	CA	980	C
1	CA	991	U
1	CA	992	U
1	CA	993	G
1	CA	1026	G
1	CA	1050	G
1	CA	1053	G
1	CA	1054	C
1	CA	1064	G
1	CA	1065	U
1	CA	1066	C
1	CA	1068	G
1	CA	1085	U
1	CA	1094	G
1	CA	1095	U
1	CA	1101	A
1	CA	1117	G
1	CA	1118	C
1	CA	1123	A
1	CA	1124	G
1	CA	1125	U
1	CA	1126	U
1	CA	1129	C
1	CA	1130	A
1	CA	1131	G
1	CA	1137	C
1	CA	1138	G
1	CA	1139	G
1	CA	1146	A

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Mol	Chain	Res	Type
1	CA	1147	C
1	CA	1152	A
1	CA	1159	U
1	CA	1184	G
1	CA	1187	G
1	CA	1190	G
1	CA	1196	U
1	CA	1197	G
1	CA	1212	U
1	CA	1213	A
1	CA	1225	A
1	CA	1227	A
1	CA	1238	A
1	CA	1249	C
1	CA	1256	A
1	CA	1257	U
1	CA	1278	U
1	CA	1280	A
1	CA	1281	U
1	CA	1282	C
1	CA	1286	A
1	CA	1287	A
1	CA	1294	G
1	CA	1297	C
1	CA	1300	G
1	CA	1301	U
1	CA	1302	U
1	CA	1305	G
1	CA	1317	C
1	CA	1320	C
1	CA	1322	C
1	CA	1325	C
1	CA	1331	G
1	CA	1335	C
1	CA	1347	G
1	CA	1364	U
1	CA	1398	A
1	CA	1419	G
1	CA	1442	G
1	CA	1442(A)	G
1	CA	1442(B)	A
1	CA	1443	G

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Mol	Chain	Res	Type
1	CA	1447	A
1	CA	1452	C
1	CA	1456	G
1	CA	1457	G
1	CA	1492	A
1	CA	1494	G
1	CA	1499	A
1	CA	1504	G
1	CA	1505	G
1	CA	1506	U
1	CA	1517	G
1	CA	1518	A
1	CA	1519	A
1	CA	1529	G
1	CA	1530	G
23	CW	5	G
23	CW	7	G
23	CW	8	U
23	CW	17	C
23	CW	18	U
23	CW	20	G
23	CW	21	U
23	CW	22	A
23	CW	24	C
23	CW	44	A
23	CW	48	U
23	CW	49	C
23	CW	54	G
23	CW	57	C
23	CW	60	A
23	CW	62	C
23	CW	66	C
23	CW	68	C
23	CW	73	A
23	CW	77	A
24	CX	13	A
24	CX	19	U
35	DA	10	G
35	DA	17	G
35	DA	28	A
35	DA	34	C
35	DA	35	G

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Mol	Chain	Res	Type
35	DA	45	C
35	DA	49	A
35	DA	50	U
35	DA	51	G
35	DA	55	G
35	DA	68	G
35	DA	69	C
35	DA	71	A
35	DA	72	U
35	DA	73	A
35	DA	75	G
35	DA	83	G
35	DA	84	A
35	DA	85	G
35	DA	88	G
35	DA	90	U
35	DA	94	C
35	DA	95	G
35	DA	99	U
35	DA	100	G
35	DA	102	G
35	DA	118	A
35	DA	120	U
35	DA	132	G
35	DA	139	G
35	DA	139(A)	G
35	DA	140	G
35	DA	142	A
35	DA	142(A)	C
35	DA	143(A)	C
35	DA	146	G
35	DA	149	A
35	DA	157	U
35	DA	158	U
35	DA	174	C
35	DA	182	A
35	DA	196	A
35	DA	197	A
35	DA	199	A
35	DA	205	G
35	DA	215	G
35	DA	216	A

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Mol	Chain	Res	Type
35	DA	221	A
35	DA	222	A
35	DA	228	A
35	DA	229	A
35	DA	248	G
35	DA	249	C
35	DA	252	G
35	DA	261	G
35	DA	271(K)	U
35	DA	271(L)	U
35	DA	271(M)	G
35	DA	271(N)	U
35	DA	271(O)	C
35	DA	271(P)	C
35	DA	271(T)	C
35	DA	271(Y)	U
35	DA	272	G
35	DA	272(B)	G
35	DA	272(H)	C
35	DA	274	G
35	DA	275	G
35	DA	283	A
35	DA	284	U
35	DA	286	C
35	DA	287	C
35	DA	292	C
35	DA	310	A
35	DA	311	A
35	DA	329	G
35	DA	330	A
35	DA	332	A
35	DA	333	G
35	DA	335	C
35	DA	343	C
35	DA	349	G
35	DA	352	G
35	DA	353	G
35	DA	356	G
35	DA	362	U
35	DA	363(F)	A
35	DA	365	C
35	DA	372	G

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Mol	Chain	Res	Type
35	DA	386	G
35	DA	387	U
35	DA	388	G
35	DA	405	U
35	DA	406	G
35	DA	411	G
35	DA	412	A
35	DA	416	C
35	DA	418	G
35	DA	428	A
35	DA	444	C
35	DA	448	U
35	DA	451	C
35	DA	456	C
35	DA	457	A
35	DA	470	A
35	DA	475	U
35	DA	481	G
35	DA	482	A
35	DA	494	G
35	DA	505	A
35	DA	508	G
35	DA	509	C
35	DA	512	G
35	DA	525	U
35	DA	529	A
35	DA	531	C
35	DA	532	A
35	DA	533	G
35	DA	537	C
35	DA	542	C
35	DA	543	C
35	DA	547	A
35	DA	548	A
35	DA	549	G
35	DA	551	G
35	DA	556	G
35	DA	562	U
35	DA	563	G
35	DA	572	A
35	DA	573	G
35	DA	575	A

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Mol	Chain	Res	Type
35	DA	586	A
35	DA	588	U
35	DA	604	G
35	DA	607	U
35	DA	614(A)	U
35	DA	614(B)	G
35	DA	614(C)	A
35	DA	615	G
35	DA	620	G
35	DA	622	G
35	DA	627	A
35	DA	637	A
35	DA	645	C
35	DA	646	A
35	DA	650	C
35	DA	652	C
35	DA	656	G
35	DA	670	A
35	DA	671	C
35	DA	686	G
35	DA	687	C
35	DA	708	C
35	DA	717	G
35	DA	722	A
35	DA	730	C
35	DA	753	C
35	DA	764	A
35	DA	765	G
35	DA	776	G
35	DA	782	A
35	DA	784	A
35	DA	785	G
35	DA	790	C
35	DA	791	C
35	DA	792	G
35	DA	805	G
35	DA	812	C
35	DA	819	A
35	DA	827	U
35	DA	828	U
35	DA	846	C
35	DA	848	G

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Mol	Chain	Res	Type
35	DA	857	C
35	DA	859	G
35	DA	864	G
35	DA	866	A
35	DA	878	A
35	DA	892	G
35	DA	896	A
35	DA	897	C
35	DA	898	C
35	DA	899	A
35	DA	904	C
35	DA	910	A
35	DA	917	A
35	DA	926	A
35	DA	932	G
35	DA	941	A
35	DA	945	A
35	DA	946	G
35	DA	955	C
35	DA	956	G
35	DA	957	A
35	DA	958	U
35	DA	959	A
35	DA	961	C
35	DA	964	C
35	DA	965	C
35	DA	973	A
35	DA	974	G
35	DA	975	C
35	DA	980	A
35	DA	983	A
35	DA	985	C
35	DA	990	A
35	DA	991	C
35	DA	996	A
35	DA	1011	G
35	DA	1012	U
35	DA	1013	C
35	DA	1017	G
35	DA	1020	A
35	DA	1022	G
35	DA	1023	U

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Mol	Chain	Res	Type
35	DA	1025	G
35	DA	1026	U
35	DA	1033	U
35	DA	1039	G
35	DA	1040	C
35	DA	1041	C
35	DA	1042	G
35	DA	1044	G
35	DA	1045	A
35	DA	1047	G
35	DA	1049	C
35	DA	1052	C
35	DA	1106	A
35	DA	1110	G
35	DA	1111	A
35	DA	1112	G
35	DA	1113	U
35	DA	1115	G
35	DA	1129	A
35	DA	1130	U
35	DA	1135	C
35	DA	1136	G
35	DA	1143	A
35	DA	1155	A
35	DA	1156	A
35	DA	1159	U
35	DA	1173	G
35	DA	1174	A
35	DA	1175	U
35	DA	1176	G
35	DA	1177	A
35	DA	1178	C
35	DA	1194	A
35	DA	1195	G
35	DA	1205	U
35	DA	1210	A
35	DA	1211	U
35	DA	1212	G
35	DA	1218	C
35	DA	1221	C
35	DA	1241	A
35	DA	1247	A

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Mol	Chain	Res	Type
35	DA	1248	G
35	DA	1249	U
35	DA	1251	C
35	DA	1253	A
35	DA	1254	A
35	DA	1255	U
35	DA	1256	G
35	DA	1265	A
35	DA	1271	G
35	DA	1272	A
35	DA	1276	A
35	DA	1281	G
35	DA	1300	U
35	DA	1301	A
35	DA	1302	A
35	DA	1314	C
35	DA	1319	G
35	DA	1329	U
35	DA	1330	C
35	DA	1332	G
35	DA	1345	C
35	DA	1349	A
35	DA	1359	A
35	DA	1368	G
35	DA	1379	A
35	DA	1380	G
35	DA	1385	G
35	DA	1386	C
35	DA	1390	U
35	DA	1407	C
35	DA	1412	A
35	DA	1416	G
35	DA	1417	C
35	DA	1420	U
35	DA	1421	G
35	DA	1428	C
35	DA	1445	A
35	DA	1449	A
35	DA	1450	G
35	DA	1460	A
35	DA	1461	G
35	DA	1467	C

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Mol	Chain	Res	Type
35	DA	1471	A
35	DA	1475	G
35	DA	1478	G
35	DA	1482	G
35	DA	1484	G
35	DA	1485	G
35	DA	1490	A
35	DA	1493	C
35	DA	1494	A
35	DA	1495	A
35	DA	1497	U
35	DA	1498	C
35	DA	1502	C
35	DA	1505	C
35	DA	1508	A
35	DA	1509	C
35	DA	1509(A)	A
35	DA	1528(A)	A
35	DA	1529	G
35	DA	1530	C
35	DA	1532	C
35	DA	1533	G
35	DA	1543	C
35	DA	1545	A
35	DA	1547	C
35	DA	1554	A
35	DA	1558	A
35	DA	1559	G
35	DA	1569	A
35	DA	1578	U
35	DA	1579	A
35	DA	1584	C
35	DA	1586	A
35	DA	1587	A
35	DA	1591	G
35	DA	1594	G
35	DA	1595	G
35	DA	1603	A
35	DA	1608	A
35	DA	1609	A
35	DA	1610	A
35	DA	1613	G

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Mol	Chain	Res	Type
35	DA	1614	A
35	DA	1615	C
35	DA	1616	A
35	DA	1617	C
35	DA	1635	G
35	DA	1640	C
35	DA	1648	C
35	DA	1654	A
35	DA	1674	G
35	DA	1678	G
35	DA	1694	C
35	DA	1696	G
35	DA	1707	G
35	DA	1718	G
35	DA	1739	U
35	DA	1741	A
35	DA	1742	G
35	DA	1746	G
35	DA	1748	G
35	DA	1763	G
35	DA	1764	G
35	DA	1773	A
35	DA	1780	A
35	DA	1791	A
35	DA	1799	G
35	DA	1800	C
35	DA	1816	G
35	DA	1820	U
35	DA	1835	G
35	DA	1847	A
35	DA	1854	A
35	DA	1858	G
35	DA	1865	G
35	DA	1866	C
35	DA	1877	A
35	DA	1878	G
35	DA	1880	C
35	DA	1882	C
35	DA	1885	A
35	DA	1888	G
35	DA	1900	A
35	DA	1903	G

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Mol	Chain	Res	Type
35	DA	1906	G
35	DA	1912	A
35	DA	1913	A
35	DA	1929	G
35	DA	1936	A
35	DA	1938	A
35	DA	1955	U
35	DA	1963	U
35	DA	1964	G
35	DA	1965	C
35	DA	1967	C
35	DA	1969	A
35	DA	1971	A
35	DA	1972	A
35	DA	1987	G
35	DA	1988	C
35	DA	1991	U
35	DA	1993	U
35	DA	2023	G
35	DA	2026	C
35	DA	2031	A
35	DA	2033	A
35	DA	2034	U
35	DA	2036	C
35	DA	2043	C
35	DA	2051	A
35	DA	2055	C
35	DA	2056	G
35	DA	2060	A
35	DA	2061	G
35	DA	2062	A
35	DA	2069	G
35	DA	2093	G
35	DA	2099	U
35	DA	2103	C
35	DA	2104	G
35	DA	2108	C
35	DA	2110	G
35	DA	2111	C
35	DA	2112	G
35	DA	2116	G
35	DA	2117	A

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Mol	Chain	Res	Type
35	DA	2118	U
35	DA	2119	A
35	DA	2120	G
35	DA	2122	U
35	DA	2127	G
35	DA	2128	C
35	DA	2163	C
35	DA	2164	C
35	DA	2165	G
35	DA	2166	G
35	DA	2169	A
35	DA	2171	A
35	DA	2172	U
35	DA	2173	A
35	DA	2176	A
35	DA	2179	C
35	DA	2185	C
35	DA	2187	G
35	DA	2190	G
35	DA	2198	A
35	DA	2199	A
35	DA	2200	C
35	DA	2207	G
35	DA	2208	A
35	DA	2218	U
35	DA	2225	A
35	DA	2226	C
35	DA	2238	G
35	DA	2239	G
35	DA	2243	U
35	DA	2251	G
35	DA	2263	C
35	DA	2275	C
35	DA	2283	C
35	DA	2285	C
35	DA	2287	A
35	DA	2290	G
35	DA	2294	C
35	DA	2297	C
35	DA	2305	A
35	DA	2307	G
35	DA	2308	G

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Mol	Chain	Res	Type
35	DA	2311	A
35	DA	2313	C
35	DA	2316	C
35	DA	2319	G
35	DA	2320	A
35	DA	2325	G
35	DA	2334	G
35	DA	2336	A
35	DA	2345	G
35	DA	2347	C
35	DA	2349	G
35	DA	2358	G
35	DA	2383	G
35	DA	2385	C
35	DA	2387	U
35	DA	2388	A
35	DA	2398	U
35	DA	2399	G
35	DA	2400	G
35	DA	2402	C
35	DA	2408	U
35	DA	2423	U
35	DA	2425	A
35	DA	2427	C
35	DA	2428	G
35	DA	2429	G
35	DA	2430	A
35	DA	2431	U
35	DA	2439	A
35	DA	2441	C
35	DA	2447	G
35	DA	2448	A
35	DA	2465	C
35	DA	2469	A
35	DA	2470	G
35	DA	2474	C
35	DA	2476	A
35	DA	2478	A
35	DA	2482	G
35	DA	2483	C
35	DA	2490	G
35	DA	2491	U

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Mol	Chain	Res	Type
35	DA	2498	C
35	DA	2502	G
35	DA	2505	G
35	DA	2518	A
35	DA	2520	C
35	DA	2523	G
35	DA	2529	G
35	DA	2534	A
35	DA	2543	G
35	DA	2554	U
35	DA	2559	C
35	DA	2566	A
35	DA	2567	G
35	DA	2571	C
35	DA	2572	A
35	DA	2573	C
35	DA	2582	G
35	DA	2585	U
35	DA	2586	C
35	DA	2602	A
35	DA	2609	U
35	DA	2610	C
35	DA	2611	U
35	DA	2612	C
35	DA	2615	U
35	DA	2630	G
35	DA	2640	G
35	DA	2646	C
35	DA	2655	G
35	DA	2673	G
35	DA	2675	A
35	DA	2682	U
35	DA	2690	C
35	DA	2691	C
35	DA	2702	U
35	DA	2703	C
35	DA	2712(A)	A
35	DA	2713	A
35	DA	2714	G
35	DA	2720	U
35	DA	2726	U
35	DA	2733	A

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Mol	Chain	Res	Type
35	DA	2752	C
35	DA	2754	U
35	DA	2759	G
35	DA	2762	G
35	DA	2763	G
35	DA	2765	A
35	DA	2766	G
35	DA	2778	A
35	DA	2779	U
35	DA	2780	G
35	DA	2781	A
35	DA	2789	C
35	DA	2790	A
35	DA	2791	C
35	DA	2796	U
35	DA	2801(A)	A
35	DA	2802	G
35	DA	2803	C
35	DA	2808	U
35	DA	2821	A
35	DA	2823	A
35	DA	2833	G
35	DA	2834	G
35	DA	2835	A
35	DA	2849	U
35	DA	2860	A
35	DA	2864	G
35	DA	2872	G
35	DA	2893	G
36	DB	8	U
36	DB	13	A
36	DB	15	A
36	DB	16	G
36	DB	22	U
36	DB	24	G
36	DB	25	A
36	DB	27	C
36	DB	33	G
36	DB	41	U
36	DB	42	C
36	DB	45	A
36	DB	47	C

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Mol	Chain	Res	Type
36	DB	53	A
36	DB	67	G
36	DB	73	A
36	DB	75	G
36	DB	88	C
36	DB	110	G

All (176) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	30	U
1	AA	60	A
1	AA	79	G
1	AA	115	G
1	AA	119	A
1	AA	243	A
1	AA	250	A
1	AA	266	G
1	AA	328	C
1	AA	366	C
1	AA	428	G
1	AA	429	U
1	AA	484	G
1	AA	509	A
1	AA	533	A
1	AA	560	U
1	AA	575	G
1	AA	687	A
1	AA	748	C
1	AA	913	A
1	AA	992	U
1	AA	1049	U
1	AA	1064	G
1	AA	1065	U
1	AA	1067	A
1	AA	1281	U
1	AA	1285	A
1	AA	1300	G
1	AA	1442(A)	G
1	AA	1498	U
1	AA	1504	G
35	BA	27	G

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Mol	Chain	Res	Type
35	BA	49	A
35	BA	71	A
35	BA	74	A
35	BA	214	G
35	BA	221	A
35	BA	272	G
35	BA	283	A
35	BA	331	A
35	BA	332	A
35	BA	334	C
35	BA	387	U
35	BA	474	G
35	BA	481	G
35	BA	542	C
35	BA	587	C
35	BA	603	A
35	BA	614(C)	A
35	BA	651	G
35	BA	669	G
35	BA	670	A
35	BA	752	A
35	BA	856	C
35	BA	1022	G
35	BA	1112	G
35	BA	1176	G
35	BA	1210	A
35	BA	1275	A
35	BA	1300	U
35	BA	1301	A
35	BA	1378	A
35	BA	1427	A
35	BA	1459	G
35	BA	1558	A
35	BA	1612	C
35	BA	1653	G
35	BA	1799	G
35	BA	1819	A
35	BA	1912	A
35	BA	1935	G
35	BA	1937	A
35	BA	1962	C
35	BA	1970	A

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Mol	Chain	Res	Type
35	BA	1987	G
35	BA	1992	G
35	BA	2033	A
35	BA	2126	A
35	BA	2128	C
35	BA	2225	A
35	BA	2282	G
35	BA	2290	G
35	BA	2313	C
35	BA	2422	A
35	BA	2481	G
35	BA	2610	C
35	BA	2689	U
35	BA	2859	G
36	BB	66	A
1	CA	30	U
1	CA	60	A
1	CA	79	G
1	CA	115	G
1	CA	119	A
1	CA	243	A
1	CA	250	A
1	CA	266	G
1	CA	328	C
1	CA	366	C
1	CA	428	G
1	CA	429	U
1	CA	484	G
1	CA	509	A
1	CA	533	A
1	CA	560	U
1	CA	575	G
1	CA	687	A
1	CA	748	C
1	CA	913	A
1	CA	992	U
1	CA	1049	U
1	CA	1064	G
1	CA	1065	U
1	CA	1067	A
1	CA	1281	U
1	CA	1285	A

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Mol	Chain	Res	Type
1	CA	1300	G
1	CA	1498	U
35	DA	27	G
35	DA	49	A
35	DA	71	A
35	DA	74	A
35	DA	214	G
35	DA	221	A
35	DA	272	G
35	DA	283	A
35	DA	331	A
35	DA	332	A
35	DA	334	C
35	DA	387	U
35	DA	474	G
35	DA	481	G
35	DA	542	C
35	DA	587	C
35	DA	603	A
35	DA	614(C)	A
35	DA	651	G
35	DA	669	G
35	DA	670	A
35	DA	752	A
35	DA	856	C
35	DA	1022	G
35	DA	1112	G
35	DA	1176	G
35	DA	1210	A
35	DA	1275	A
35	DA	1300	U
35	DA	1301	A
35	DA	1378	A
35	DA	1427	A
35	DA	1459	G
35	DA	1558	A
35	DA	1612	C
35	DA	1653	G
35	DA	1799	G
35	DA	1819	A
35	DA	1912	A
35	DA	1935	G

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Mol	Chain	Res	Type
35	DA	1937	A
35	DA	1962	C
35	DA	1970	A
35	DA	1987	G
35	DA	1992	G
35	DA	2033	A
35	DA	2126	A
35	DA	2128	C
35	DA	2225	A
35	DA	2282	G
35	DA	2313	C
35	DA	2422	A
35	DA	2481	G
35	DA	2610	C
35	DA	2689	U
35	DA	2763	G
35	DA	2859	G
36	DB	66	A

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
23	5MU	CW	55	23	15,22,23	1.22	2 (13%)	16,32,35	3.72	1 (6%)
23	5MU	AW	55	23	15,22,23	1.18	2 (13%)	16,32,35	3.77	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
23	5MU	CW	55	23	-	0/5/25/26	0/2/2/2
23	5MU	AW	55	23	-	0/5/25/26	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	CW	55	5MU	C4-N3	3.42	1.39	1.33
23	AW	55	5MU	C4-N3	3.41	1.39	1.33
23	CW	55	5MU	C6-C5	-2.13	1.34	1.40
23	AW	55	5MU	C6-C5	-2.01	1.34	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	AW	55	5MU	C4-N3-C2	14.73	127.57	115.14
23	CW	55	5MU	C4-N3-C2	14.48	127.36	115.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
23	CW	55	5MU	3	0
23	AW	55	5MU	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1486 ligands modelled in this entry, 1486 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
13	CM	5
13	AM	5
9	AI	2
9	CI	2
52	DV	1
41	DG	1
41	BG	1
52	BV	1
32	D6	1
32	B6	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B6	46:HIS	C	47:THR	N	6.15
1	D6	46:HIS	C	47:THR	N	6.12
1	CI	53:VAL	C	54:ASP	N	5.30
1	AI	53:VAL	C	54:ASP	N	5.29
1	DG	112:PRO	C	113:ARG	N	4.36
1	CM	69:GLU	C	70:LEU	N	4.32
1	AM	69:GLU	C	70:LEU	N	4.30
1	DV	80:GLN	C	81:TYR	N	3.61
1	CM	112:GLY	C	113:PRO	N	3.58
1	AM	112:GLY	C	113:PRO	N	3.57
1	BV	80:GLN	C	81:TYR	N	3.57
1	CM	118:ALA	C	119:GLY	N	3.52
1	AM	118:ALA	C	119:GLY	N	3.48
1	CI	104:ARG	C	105:ASP	N	3.03
1	BG	112:PRO	C	113:ARG	N	3.00
1	AI	104:ARG	C	105:ASP	N	2.98
1	AM	97:PRO	C	98:VAL	N	2.93
1	CM	97:PRO	C	98:VAL	N	2.92

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	CM	65:LYS	C	66:LEU	N	2.52
1	AM	65:LYS	C	66:LEU	N	2.50

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AA	1504/1522 (98%)	0.64	162 (10%) 5 6	15, 82, 177, 200	0
1	CA	1504/1522 (98%)	0.81	195 (12%) 3 4	20, 90, 190, 200	0
2	AB	235/256 (91%)	0.21	22 (9%) 8 9	19, 113, 188, 200	0
2	CB	235/256 (91%)	0.22	21 (8%) 9 10	26, 127, 188, 200	0
3	AC	207/239 (86%)	0.11	14 (6%) 17 16	11, 98, 175, 200	0
3	CC	207/239 (86%)	0.27	16 (7%) 13 13	38, 121, 183, 200	0
4	AD	208/209 (99%)	-0.23	4 (1%) 66 61	13, 75, 145, 200	0
4	CD	208/209 (99%)	-0.23	3 (1%) 75 69	25, 89, 155, 200	0
5	AE	151/162 (93%)	-0.22	3 (1%) 65 60	21, 73, 158, 190	0
5	CE	151/162 (93%)	0.11	6 (3%) 38 33	26, 89, 157, 200	0
6	AF	101/101 (100%)	-0.15	2 (1%) 65 60	19, 100, 151, 189	0
6	CF	101/101 (100%)	-0.24	0 100 100	16, 83, 141, 170	0
7	AG	155/156 (99%)	0.16	16 (10%) 6 7	34, 104, 175, 194	0
7	CG	155/156 (99%)	0.31	18 (11%) 4 5	53, 123, 185, 200	0
8	AH	138/138 (100%)	-0.31	1 (0%) 87 83	33, 81, 145, 190	0
8	CH	138/138 (100%)	-0.37	1 (0%) 87 83	30, 91, 147, 165	0
9	AI	127/128 (99%)	0.00	4 (3%) 49 43	27, 123, 180, 200	0
9	CI	127/128 (99%)	0.80	27 (21%) 0 1	43, 140, 195, 200	0
10	AJ	99/105 (94%)	0.23	5 (5%) 28 25	47, 120, 180, 200	0
10	CJ	99/105 (94%)	1.01	16 (16%) 1 2	65, 140, 189, 200	0
11	AK	119/129 (92%)	0.27	12 (10%) 7 7	13, 78, 163, 187	0
11	CK	119/129 (92%)	0.30	9 (7%) 13 14	30, 86, 164, 200	0
12	AL	125/135 (92%)	-0.36	1 (0%) 86 81	19, 71, 146, 200	0
12	CL	125/135 (92%)	-0.26	1 (0%) 86 81	12, 75, 143, 200	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AM	125/126 (99%)	0.55	20 (16%) 1 2	56, 124, 183, 200	0
13	CM	125/126 (99%)	0.32	12 (9%) 8 8	41, 140, 199, 200	0
14	AN	60/61 (98%)	-0.12	2 (3%) 46 41	38, 94, 181, 200	0
14	CN	60/61 (98%)	0.06	1 (1%) 70 64	57, 117, 171, 189	0
15	AO	88/89 (98%)	-0.37	1 (1%) 80 75	11, 84, 147, 172	0
15	CO	88/89 (98%)	-0.18	4 (4%) 33 29	18, 88, 146, 167	0
16	AP	84/88 (95%)	-0.23	2 (2%) 59 53	17, 64, 135, 200	0
16	CP	84/88 (95%)	0.21	8 (9%) 8 8	32, 90, 139, 199	0
17	AQ	100/105 (95%)	0.11	11 (11%) 5 6	15, 81, 143, 200	0
17	CQ	100/105 (95%)	0.23	14 (14%) 2 3	20, 84, 134, 200	0
18	AR	70/88 (79%)	0.09	2 (2%) 51 45	33, 87, 156, 200	0
18	CR	70/88 (79%)	0.00	4 (5%) 23 21	17, 79, 137, 170	0
19	AS	79/93 (84%)	0.14	3 (3%) 40 36	60, 134, 200, 200	0
19	CS	79/93 (84%)	0.47	11 (13%) 2 3	76, 140, 200, 200	0
20	AT	99/106 (93%)	0.11	6 (6%) 21 19	21, 91, 178, 200	0
20	CT	99/106 (93%)	-0.03	7 (7%) 16 15	45, 102, 167, 197	0
21	AU	25/27 (92%)	0.06	0 100 100	59, 114, 161, 200	0
21	CU	25/27 (92%)	0.98	5 (20%) 1 1	58, 131, 182, 200	0
22	AV	17/76 (22%)	0.68	2 (11%) 4 5	55, 68, 122, 125	0
22	CV	17/76 (22%)	0.33	1 (5%) 22 20	56, 73, 128, 131	0
23	AW	76/77 (98%)	3.55	45 (59%) 0 0	63, 179, 200, 200	0
23	CW	76/77 (98%)	5.00	61 (80%) 0 0	62, 191, 200, 200	0
24	AX	11/31 (35%)	-0.04	0 100 100	23, 72, 120, 172	0
24	CX	11/31 (35%)	-0.05	0 100 100	41, 63, 93, 196	0
25	AY	185/185 (100%)	0.60	30 (16%) 1 2	27, 99, 167, 200	0
25	CY	185/185 (100%)	0.38	27 (14%) 2 3	9, 93, 166, 198	0
26	B0	85/85 (100%)	0.59	11 (12%) 3 4	21, 80, 184, 199	0
26	D0	85/85 (100%)	0.64	11 (12%) 3 4	39, 95, 174, 200	0
27	B1	89/98 (90%)	-0.30	2 (2%) 62 56	5, 64, 138, 193	0
27	D1	89/98 (90%)	-0.49	2 (2%) 62 56	5, 62, 133, 200	0
28	B2	51/72 (70%)	-0.31	2 (3%) 39 35	44, 103, 157, 200	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	D2	51/72 (70%)	0.13	5 (9%) 7 8	24, 93, 168, 200	0
29	B3	60/60 (100%)	-0.13	1 (1%) 70 64	15, 81, 158, 200	0
29	D3	60/60 (100%)	-0.14	1 (1%) 70 64	21, 85, 167, 200	0
30	B4	50/71 (70%)	-0.18	0 100 100	71, 144, 192, 200	0
30	D4	50/71 (70%)	0.18	5 (10%) 7 8	86, 167, 192, 200	0
31	B5	59/60 (98%)	0.09	4 (6%) 17 16	22, 84, 184, 200	0
31	D5	59/60 (98%)	-0.35	0 100 100	6, 74, 191, 200	0
32	B6	45/54 (83%)	0.83	7 (15%) 2 2	53, 122, 184, 200	0
32	D6	45/54 (83%)	1.44	14 (31%) 0 0	54, 147, 188, 200	0
33	B7	49/49 (100%)	-0.23	0 100 100	5, 47, 114, 200	0
33	D7	49/49 (100%)	-0.38	1 (2%) 65 60	3, 44, 105, 176	0
34	B8	64/65 (98%)	-0.17	1 (1%) 72 66	7, 53, 145, 200	0
34	D8	64/65 (98%)	-0.12	2 (3%) 49 43	27, 73, 140, 185	0
35	BA	2767/2782 (99%)	0.70	328 (11%) 4 5	9, 65, 172, 200	0
35	DA	2767/2782 (99%)	0.70	321 (11%) 4 5	6, 62, 169, 200	0
36	BB	119/122 (97%)	1.06	17 (14%) 2 3	60, 102, 160, 189	0
36	DB	119/122 (97%)	1.13	25 (21%) 1 1	69, 140, 188, 200	0
37	BC	191/229 (83%)	2.09	75 (39%) 0 0	76, 175, 200, 200	0
37	DC	191/229 (83%)	2.60	99 (51%) 0 0	120, 180, 200, 200	0
38	BD	272/276 (98%)	-0.40	1 (0%) 92 90	8, 56, 122, 200	0
38	DD	272/276 (98%)	-0.47	2 (0%) 87 83	1, 47, 114, 200	0
39	BE	205/206 (99%)	0.12	11 (5%) 25 23	10, 75, 160, 200	0
39	DE	205/206 (99%)	-0.14	5 (2%) 59 53	7, 64, 153, 199	0
40	BF	208/210 (99%)	-0.08	10 (4%) 30 27	1, 69, 164, 200	0
40	DF	208/210 (99%)	-0.15	6 (2%) 51 45	6, 75, 170, 200	0
41	BG	181/182 (99%)	0.29	22 (12%) 4 5	36, 107, 171, 200	0
41	DG	181/182 (99%)	0.78	28 (15%) 2 2	45, 135, 191, 200	0
42	BH	160/180 (88%)	0.62	23 (14%) 2 3	47, 145, 200, 200	0
42	DH	160/180 (88%)	0.09	9 (5%) 24 22	36, 110, 180, 200	0
43	BI	146/148 (98%)	0.29	15 (10%) 6 7	25, 112, 184, 200	0
43	DI	146/148 (98%)	1.75	45 (30%) 0 0	17, 134, 200, 200	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	BN	139/140 (99%)	-0.11	7 (5%) 28 25	37, 98, 160, 200	0
44	DN	139/140 (99%)	-0.31	4 (2%) 51 45	19, 76, 152, 195	0
45	BO	122/122 (100%)	-0.25	3 (2%) 57 51	15, 63, 116, 133	0
45	DO	122/122 (100%)	-0.41	0 100 100	10, 49, 100, 128	0
46	BP	146/150 (97%)	0.06	10 (6%) 17 16	6, 84, 167, 200	0
46	DP	146/150 (97%)	0.13	8 (5%) 25 22	18, 93, 168, 199	0
47	BQ	136/141 (96%)	-0.00	7 (5%) 28 25	26, 90, 187, 200	0
47	DQ	136/141 (96%)	-0.25	3 (2%) 62 56	16, 73, 158, 200	0
48	BR	117/118 (99%)	-0.47	0 100 100	8, 66, 138, 198	0
48	DR	117/118 (99%)	-0.43	0 100 100	8, 64, 136, 180	0
49	BS	99/112 (88%)	0.22	7 (7%) 16 15	32, 105, 186, 200	0
49	DS	99/112 (88%)	1.01	23 (23%) 0 0	62, 133, 182, 200	0
50	BT	138/146 (94%)	-0.12	4 (2%) 51 45	21, 96, 190, 200	0
50	DT	138/146 (94%)	-0.07	8 (5%) 23 20	17, 84, 173, 200	0
51	BU	117/118 (99%)	-0.25	4 (3%) 45 40	15, 79, 161, 182	0
51	DU	117/118 (99%)	-0.20	5 (4%) 35 31	12, 65, 141, 194	0
52	BV	101/101 (100%)	0.57	16 (15%) 2 2	26, 106, 188, 200	0
52	DV	101/101 (100%)	0.09	9 (8%) 9 10	29, 100, 178, 200	0
53	BW	113/113 (100%)	-0.45	2 (1%) 68 62	8, 55, 134, 200	0
53	DW	113/113 (100%)	-0.34	2 (1%) 68 62	16, 66, 148, 200	0
54	BX	93/96 (96%)	-0.35	1 (1%) 80 75	12, 83, 163, 200	0
54	DX	93/96 (96%)	-0.35	0 100 100	18, 79, 147, 185	0
55	BY	101/110 (91%)	0.03	8 (7%) 12 13	16, 94, 187, 200	0
55	DY	101/110 (91%)	0.35	11 (10%) 5 6	21, 108, 187, 200	0
56	BZ	177/206 (85%)	-0.12	7 (3%) 38 33	46, 113, 178, 200	0
56	DZ	177/206 (85%)	0.22	13 (7%) 15 15	11, 114, 180, 200	0
All	All	21176/22108 (95%)	0.41	2111 (9%) 7 8	1, 84, 184, 200	0

All (2111) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
35	BA	897	C	19.0
35	DA	2109	U	18.4

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Mol	Chain	Res	Type	RSRZ
35	DA	897	C	18.4
43	DI	143	SER	17.7
23	CW	7	G	14.9
35	DA	2174	C	13.6
55	DY	52	SER	13.5
37	DC	179	SER	12.9
1	CA	88	A	12.9
37	BC	178	ALA	12.6
23	CW	35	C	12.6
32	D6	43	CYS	12.5
37	DC	38	ASP	12.3
43	DI	67	ARG	12.3
37	BC	90	GLY	12.3
43	DI	66	GLU	12.2
1	CA	89	C	12.2
23	CW	50	G	12.0
37	DC	181	PRO	11.7
43	DI	70	GLU	11.6
42	BH	44	VAL	11.4
23	CW	6	G	11.4
37	BC	77	ILE	11.3
41	DG	134	GLY	11.3
1	CA	1027	C	11.3
35	DA	1531	C	11.2
35	DA	2110	G	11.1
37	BC	40	THR	11.0
1	CA	1009	G	10.8
1	AA	1027	C	10.8
35	DA	2168	G	10.7
41	DG	89	GLY	10.3
23	AW	16	C	10.1
23	AW	7	G	10.1
43	DI	92	VAL	10.0
10	CJ	71	LEU	9.8
37	BC	34	THR	9.8
1	AA	1026	G	9.7
37	BC	91	ALA	9.7
1	CA	81	U	9.7
43	DI	71	ILE	9.6
41	DG	135	LEU	9.5
23	CW	11	A	9.5
43	DI	90	GLY	9.5

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Mol	Chain	Res	Type	RSRZ
10	CJ	38	ILE	9.4
35	DA	2120	G	9.4
23	AW	50	G	9.3
23	AW	60	A	9.3
1	CA	80	G	9.3
43	DI	108	THR	9.2
43	DI	119	PRO	9.2
37	DC	180	PHE	9.2
23	CW	33	C	9.2
1	CA	1026	G	9.1
41	DG	88	ILE	9.0
35	DA	2802	G	9.0
35	BA	2792	G	9.0
26	D0	6	ALA	8.9
23	CW	57	C	8.9
35	BA	1048	A	8.9
35	BA	2169	A	8.9
35	BA	334	C	8.9
23	CW	2	G	8.8
23	CW	36	A	8.8
26	D0	7	LEU	8.8
35	DA	2108	C	8.8
43	DI	84	GLY	8.6
27	D1	96	LYS	8.6
27	B1	96	LYS	8.6
35	DA	2169	A	8.6
1	AA	1034	G	8.6
37	DC	183	GLU	8.5
23	CW	24	C	8.5
41	DG	41	GLN	8.5
47	BQ	140	ALA	8.4
1	AA	76	C	8.4
1	AA	1030(B)	C	8.3
35	DA	2792	G	8.3
35	BA	2791	C	8.2
35	DA	2892	A	8.2
35	BA	2892	A	8.2
35	BA	2893	G	8.2
23	AW	15	G	8.1
35	BA	2793	G	8.1
23	CW	58	A	8.1
37	DC	78	ALA	8.1

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Mol	Chain	Res	Type	RSRZ
23	CW	62	C	8.1
23	CW	54	G	7.9
37	DC	122	ALA	7.9
23	CW	34	U	7.8
26	B0	6	ALA	7.8
37	DC	172	HIS	7.8
1	CA	82	U	7.8
23	CW	56	U	7.8
26	B0	3	HIS	7.7
55	DY	50	ARG	7.7
25	CY	70	SER	7.7
46	DP	119	GLU	7.7
37	DC	222	VAL	7.7
23	AW	8	U	7.6
23	AW	52	C	7.6
31	B5	60	VAL	7.6
37	DC	85	GLU	7.6
37	BC	39	GLU	7.5
23	CW	32	G	7.5
23	CW	8	U	7.5
1	CA	90	U	7.5
35	DA	2175	C	7.5
35	BA	2115	G	7.4
37	BC	95	GLY	7.4
50	DT	2	ASN	7.4
23	CW	10	G	7.4
26	B0	5	LYS	7.4
35	DA	2126	A	7.4
1	AA	73	G	7.3
23	AW	35	C	7.3
35	DA	895	U	7.3
23	AW	61	U	7.3
35	BA	1496	A	7.3
37	DC	178	ALA	7.3
35	BA	899	A	7.3
37	DC	46	LYS	7.2
17	CQ	69	LYS	7.2
37	BC	41	VAL	7.2
41	DG	155	MET	7.2
55	BY	61	ILE	7.2
47	BQ	141	GLN	7.1
23	AW	53	G	7.1

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Mol	Chain	Res	Type	RSRZ
41	BG	89	GLY	7.1
37	DC	182	PRO	7.1
23	AW	13	C	7.1
35	BA	2104	G	7.1
35	DA	2113	U	7.1
35	DA	2790	A	7.0
35	DA	2115	G	7.0
23	CW	22	A	7.0
37	DC	176	GLY	7.0
32	B6	45	LYS	7.0
35	DA	2116	G	7.0
37	DC	23	ASP	7.0
56	DZ	27	VAL	6.9
37	BC	121	GLY	6.8
50	DT	138	ALA	6.8
23	CW	5	G	6.8
37	BC	120	MET	6.8
43	DI	53	ALA	6.8
23	AW	20	G	6.8
37	BC	124	GLY	6.8
37	DC	88	GLU	6.7
37	BC	78	ALA	6.7
35	DA	2117	A	6.7
41	DG	49	ASP	6.7
41	BG	42	GLY	6.7
1	AA	97	G	6.7
35	BA	2168	G	6.7
43	DI	69	LYS	6.6
23	CW	49	C	6.6
35	DA	894	C	6.6
13	AM	124	PRO	6.6
23	CW	23	G	6.6
23	CW	9	G	6.6
7	AG	84	ASN	6.6
1	AA	1005	A	6.6
23	AW	58	A	6.6
35	DA	2114	A	6.6
35	DA	884	C	6.6
35	DA	2170	A	6.6
35	BA	2110	G	6.5
43	DI	141	LYS	6.5
2	AB	228	GLY	6.5

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Mol	Chain	Res	Type	RSRZ
37	BC	119	VAL	6.5
37	BC	96	GLY	6.5
35	BA	1110	G	6.5
43	DI	125	GLU	6.5
32	D6	42	TRP	6.5
37	BC	88	GLU	6.5
1	CA	993	G	6.5
35	DA	1448	G	6.5
37	DC	37	PHE	6.5
37	DC	96	GLY	6.4
47	BQ	139	GLU	6.4
36	DB	4	C	6.4
35	DA	2173	A	6.4
35	BA	1527	G	6.4
37	BC	155	GLU	6.4
35	BA	2170	A	6.3
35	BA	279	C	6.3
1	CA	79	G	6.3
35	BA	1450	G	6.3
42	DH	45	VAL	6.2
35	BA	2173	A	6.2
10	CJ	83	GLU	6.2
2	AB	35	GLU	6.2
1	AA	1003	G	6.2
1	CA	153	C	6.2
23	CW	63	C	6.2
23	AW	59	A	6.2
42	BH	24	VAL	6.2
43	DI	97	ILE	6.2
1	CA	83	U	6.2
3	AC	102	ASN	6.1
37	DC	187	ASP	6.1
41	BG	39	ILE	6.1
35	BA	1507	A	6.1
23	AW	9	G	6.1
23	CW	61	U	6.0
1	CA	1028	C	6.0
1	CA	1043	C	6.0
37	BC	35	ALA	6.0
41	DG	40	ASN	6.0
55	BY	50	ARG	6.0
35	DA	2893	G	6.0

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Mol	Chain	Res	Type	RSRZ
23	AW	22	A	6.0
23	CW	26	C	6.0
23	CW	51	U	5.9
23	AW	33	C	5.9
35	DA	271(S)	G	5.9
35	DA	2801	A	5.9
23	AW	34	U	5.9
35	DA	75	G	5.9
35	DA	2789	C	5.9
7	CG	156	TRP	5.9
19	CS	82	GLY	5.9
47	DQ	140	ALA	5.9
35	BA	1578	U	5.9
37	BC	122	ALA	5.9
26	D0	8	ALA	5.9
37	DC	81	GLU	5.8
1	CA	1025	U	5.8
35	BA	883	G	5.8
35	DA	275	G	5.8
35	BA	2189	U	5.8
32	D6	40	CYS	5.8
35	DA	1044	G	5.8
23	CW	72	C	5.8
23	AW	23	G	5.8
13	CM	24	GLY	5.8
40	BF	12	LEU	5.7
35	DA	652	C	5.7
26	B0	4	LYS	5.7
1	CA	1141	C	5.7
35	BA	2114	A	5.7
35	BA	2167	U	5.7
1	AA	80	G	5.7
35	DA	272(J)	C	5.7
35	DA	1508	A	5.6
50	BT	135	ALA	5.6
37	DC	82	LYS	5.6
1	AA	1002	G	5.6
9	AI	18	PHE	5.6
23	CW	12	G	5.6
23	CW	46	G	5.6
35	DA	271(R)	G	5.6
35	DA	2121	G	5.6

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Mol	Chain	Res	Type	RSRZ
55	DY	59	GLY	5.6
10	CJ	73	ASP	5.6
35	BA	2125	G	5.6
1	AA	1035	A	5.6
35	DA	2793	G	5.6
23	CW	4	G	5.5
46	BP	149	GLU	5.5
43	BI	146	ALA	5.5
37	BC	125	SER	5.5
1	CA	425	G	5.5
35	BA	1506	C	5.5
20	CT	104	LEU	5.5
1	CA	84	U	5.5
37	DC	21	THR	5.5
1	CA	1031	G	5.5
35	BA	1509(A)	A	5.5
23	AW	6	G	5.4
55	BY	60	PHE	5.4
1	CA	1129	C	5.4
35	BA	414	C	5.4
1	AA	993	G	5.4
35	DA	1505	C	5.4
1	CA	1030(C)	G	5.4
13	AM	125	ARG	5.4
13	AM	84	ILE	5.4
35	DA	2124	G	5.4
1	CA	1030(B)	C	5.4
43	DI	64	GLU	5.4
37	BC	179	SER	5.3
1	CA	1295	G	5.3
23	AW	24	C	5.3
47	DQ	141	GLN	5.3
13	AM	7	VAL	5.3
15	CO	23	GLY	5.3
56	DZ	28	MET	5.3
35	BA	2179	C	5.3
35	BA	2166	G	5.3
37	DC	190	ARG	5.3
37	DC	123	VAL	5.3
35	BA	1865	G	5.3
35	DA	2791	C	5.3
35	BA	1044	G	5.2

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Mol	Chain	Res	Type	RSRZ
43	DI	107	VAL	5.2
35	DA	1865	G	5.2
23	AW	49	C	5.2
23	CW	15	G	5.2
23	CW	25	U	5.2
37	DC	45	ALA	5.2
35	BA	1467	C	5.2
36	BB	5	C	5.2
41	BG	43	LEU	5.2
3	AC	82	GLU	5.2
9	CI	96	LEU	5.2
37	DC	97	GLU	5.2
1	CA	1034	G	5.2
35	BA	1422	G	5.2
7	CG	53	LYS	5.1
19	CS	40	ILE	5.1
25	CY	59	THR	5.1
37	DC	216	THR	5.1
35	BA	2808	U	5.1
1	AA	1445	C	5.1
2	CB	227	GLY	5.1
37	DC	95	GLY	5.1
43	DI	130	TYR	5.1
1	CA	1024	G	5.1
1	AA	1260	C	5.1
25	CY	102	ASN	5.1
36	DB	49	C	5.1
25	AY	52	LEU	5.1
1	AA	1021	G	5.1
35	BA	1017	G	5.1
10	CJ	4	ILE	5.1
37	BC	108	MET	5.1
55	DY	53	PRO	5.1
35	DA	1544	A	5.1
32	D6	39	TYR	5.0
35	BA	1531	C	5.0
37	BC	38	ASP	5.0
23	CW	14	A	5.0
35	BA	1174	A	5.0
35	BA	356	G	5.0
41	DG	90	LEU	5.0
52	DV	46	VAL	5.0

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Mol	Chain	Res	Type	RSRZ
35	DA	2105	C	5.0
35	DA	2167	U	5.0
43	DI	120	ILE	5.0
35	DA	1530	C	5.0
35	DA	290	G	5.0
1	AA	1220	G	5.0
35	DA	11	G	5.0
26	D0	85	ALA	5.0
23	AW	25	U	5.0
37	DC	215	THR	5.0
35	DA	2127	G	5.0
23	CW	37	U	5.0
35	DA	2125	G	5.0
36	BB	107	G	5.0
23	AW	51	U	4.9
3	AC	66	VAL	4.9
25	CY	100	TYR	4.9
23	CW	47	G	4.9
35	DA	604	G	4.9
42	BH	19	VAL	4.9
1	CA	1005	A	4.9
25	CY	67	VAL	4.9
23	AW	47	G	4.9
2	AB	40	HIS	4.9
9	CI	22	GLY	4.9
35	DA	2123	G	4.9
2	CB	207	ALA	4.9
56	DZ	173	ALA	4.9
50	DT	39	ARG	4.9
35	BA	2467	C	4.9
3	CC	143	GLU	4.9
35	DA	1048	A	4.9
52	BV	67	GLY	4.9
37	DC	43	VAL	4.9
35	DA	2891	G	4.9
43	DI	94	ALA	4.9
37	DC	94	VAL	4.9
1	CA	97	G	4.9
11	CK	80	VAL	4.9
37	BC	24	GLU	4.9
1	AA	83	U	4.8
35	DA	877	U	4.8

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Mol	Chain	Res	Type	RSRZ
7	AG	86	GLN	4.8
35	DA	2468	G	4.8
35	DA	312	G	4.8
1	CA	1163	C	4.8
23	AW	66	C	4.8
35	BA	301	G	4.8
23	CW	21	U	4.8
23	AW	12	G	4.8
35	BA	652	C	4.8
35	BA	1526	G	4.8
43	DI	127	VAL	4.8
35	BA	1016	G	4.8
1	AA	89	C	4.8
35	BA	2477	C	4.8
1	CA	688	G	4.8
25	CY	98	ALA	4.8
2	CB	128	GLU	4.8
37	BC	87	GLU	4.8
25	AY	97	ASP	4.7
35	DA	1527	G	4.7
35	DA	2477	C	4.7
35	BA	2113	U	4.7
37	DC	89	ALA	4.7
35	DA	2186	G	4.7
37	BC	123	VAL	4.7
10	CJ	5	ARG	4.7
37	DC	221	SER	4.7
1	CA	1128	C	4.7
42	BH	46	GLU	4.7
49	DS	55	ALA	4.7
21	CU	26	LYS	4.7
3	AC	79	ARG	4.7
1	CA	70	G	4.7
35	BA	357	A	4.7
25	CY	73	GLN	4.7
35	DA	1174	A	4.6
23	AW	11	A	4.6
35	BA	1508	A	4.6
35	DA	1045	A	4.6
35	BA	1115	G	4.6
37	BC	173	ALA	4.6
36	BB	66	A	4.6

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Mol	Chain	Res	Type	RSRZ
35	BA	2124	G	4.6
41	BG	127	GLY	4.6
1	AA	992	U	4.6
13	AM	119	GLY	4.6
37	DC	18	LYS	4.6
1	CA	218	C	4.6
7	AG	85	TYR	4.6
37	DC	208	PHE	4.6
37	BC	60	GLY	4.6
41	DG	87	PRO	4.6
35	BA	2796	U	4.6
17	AQ	16	GLN	4.5
35	BA	2801	A	4.5
23	CW	53	G	4.5
23	AW	57	C	4.5
1	AA	1006	C	4.5
1	CA	71	C	4.5
23	AW	32	G	4.5
23	CW	1	C	4.5
35	BA	2471	C	4.5
35	DA	352	G	4.5
26	B0	2	ALA	4.5
37	BC	83	ILE	4.5
40	BF	133	ASN	4.5
23	CW	59	A	4.5
35	BA	2476	A	4.5
35	BA	1577	C	4.5
1	AA	1086	U	4.5
15	CO	22	THR	4.5
10	CJ	37	PRO	4.5
42	DH	82	GLY	4.5
1	CA	1217	C	4.5
35	BA	1714	G	4.5
35	BA	2532	G	4.5
17	CQ	64	PRO	4.5
37	BC	57	ASN	4.5
23	CW	45	A	4.5
30	D4	1	MET	4.5
2	AB	90	MET	4.5
35	DA	2665	A	4.5
35	BA	893	C	4.5
35	BA	355	G	4.4

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Mol	Chain	Res	Type	RSRZ
35	BA	2120	G	4.4
37	DC	39	GLU	4.4
35	BA	2165	G	4.4
7	CG	74	GLU	4.4
17	AQ	69	LYS	4.4
1	AA	79	G	4.4
1	CA	922	G	4.4
1	CA	1021	G	4.4
20	AT	104	LEU	4.4
35	BA	2121	G	4.4
35	BA	2126	A	4.4
35	DA	896	A	4.4
35	BA	2105	C	4.4
17	AQ	18	THR	4.4
2	CB	18	GLY	4.4
41	BG	38	VAL	4.4
1	CA	152	A	4.4
35	BA	401	A	4.4
35	BA	884	C	4.4
15	CO	24	SER	4.4
26	B0	7	LEU	4.4
40	BF	24	LEU	4.4
37	DC	74	VAL	4.4
43	BI	43	ASN	4.4
40	DF	23	ASP	4.4
19	AS	82	GLY	4.4
35	DA	1863	G	4.4
36	BB	106	G	4.4
35	DA	2476	A	4.3
40	BF	134	GLY	4.3
1	CA	570	G	4.3
3	AC	101	LEU	4.3
43	DI	144	VAL	4.3
1	AA	82	U	4.3
1	CA	1238	A	4.3
7	CG	4	ARG	4.3
37	BC	211	SER	4.3
18	AR	77	GLY	4.3
35	BA	2106	G	4.3
35	BA	2567	G	4.3
35	BA	2795	G	4.3
23	CW	52	C	4.3

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Mol	Chain	Res	Type	RSRZ
35	BA	2108	C	4.3
25	AY	99	LEU	4.3
35	BA	2112	G	4.3
35	DA	74	A	4.3
37	DC	133	PRO	4.3
37	DC	90	GLY	4.3
41	BG	49	ASP	4.3
9	CI	95	LYS	4.3
35	BA	708	C	4.3
37	DC	132	ALA	4.3
35	BA	1111	A	4.3
20	AT	103	GLY	4.3
22	AV	26	G	4.3
35	DA	2795	G	4.3
40	DF	133	ASN	4.3
36	BB	4	C	4.3
1	AA	1025	U	4.2
9	CI	33	PHE	4.2
35	BA	1176	G	4.2
35	BA	2116	G	4.2
35	BA	2410	G	4.2
35	DA	2166	G	4.2
10	CJ	39	PRO	4.2
37	DC	154	ARG	4.2
25	CY	74	ASN	4.2
43	BI	58	LEU	4.2
1	AA	1259	C	4.2
1	AA	1008	C	4.2
42	BH	43	VAL	4.2
25	AY	55	ILE	4.2
51	DU	75	ASN	4.2
35	BA	2164	C	4.2
35	DA	2799	C	4.2
2	AB	208	ILE	4.2
39	BE	33	VAL	4.2
25	AY	100	TYR	4.2
35	BA	290	G	4.2
35	DA	271(Q)	G	4.2
35	DA	1465	G	4.2
35	BA	1029	A	4.2
2	AB	41	ILE	4.2
32	B6	14	THR	4.2

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Mol	Chain	Res	Type	RSRZ
37	BC	80	GLY	4.2
37	DC	211	SER	4.2
35	BA	289	A	4.2
30	D4	2	LYS	4.2
42	DH	46	GLU	4.2
49	DS	33	LYS	4.2
18	CR	51	LEU	4.1
35	BA	2894	G	4.1
35	BA	1505	C	4.1
41	DG	156	ASP	4.1
1	CA	1033	G	4.1
35	BA	94(A)	G	4.1
35	DA	226	G	4.1
43	DI	139	GLN	4.1
37	BC	79	LYS	4.1
37	DC	157	LYS	4.1
1	AA	1033	G	4.1
1	AA	1029	C	4.1
35	BA	1546	C	4.1
25	AY	47	GLY	4.1
35	BA	646	A	4.1
35	DA	2808	U	4.1
25	CY	101	ILE	4.1
25	CY	58	VAL	4.1
43	DI	110	ASP	4.1
1	AA	1129	C	4.1
35	DA	271(I)	G	4.1
35	DA	1466	G	4.1
55	BY	52	SER	4.1
35	DA	2467	C	4.1
37	DC	158	ALA	4.1
23	CW	27	G	4.1
35	BA	30	G	4.1
1	CA	98	G	4.1
1	CA	1083	U	4.1
23	CW	20	G	4.1
23	AW	26	C	4.1
23	AW	67	C	4.1
23	CW	64	G	4.1
35	BA	2662	A	4.1
25	CY	72	ASP	4.1
35	BA	2473	U	4.0

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Mol	Chain	Res	Type	RSRZ
1	CA	838	G	4.0
1	CA	1003	G	4.0
35	BA	1215	G	4.0
35	DA	1171	G	4.0
35	DA	1447	G	4.0
35	DA	1450	G	4.0
28	D2	43	GLN	4.0
52	BV	45	THR	4.0
35	BA	2637	U	4.0
35	BA	2799	C	4.0
35	BA	415	A	4.0
35	DA	899	A	4.0
56	DZ	88	PHE	4.0
35	BA	2857	G	4.0
37	DC	93	TYR	4.0
31	B5	53	ALA	4.0
1	CA	689	C	4.0
13	AM	123	ALA	4.0
37	DC	152	ILE	4.0
35	BA	2122	U	4.0
35	DA	1436	G	4.0
40	BF	25	PRO	4.0
1	CA	1094	G	4.0
4	AD	36	ARG	4.0
25	CY	97	ASP	4.0
37	BC	23	ASP	4.0
1	CA	426	G	4.0
23	AW	14	A	4.0
35	DA	2809	A	4.0
52	BV	68	LYS	4.0
37	BC	172	HIS	4.0
43	DI	63	ALA	4.0
43	DI	85	GLU	4.0
35	BA	898	C	4.0
35	DA	645	C	4.0
35	DA	1422	G	4.0
37	BC	109	ASP	4.0
35	DA	1864	U	4.0
47	BQ	123	HIS	4.0
23	CW	48	U	3.9
35	BA	2470	G	3.9
1	CA	1502	A	3.9

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Mol	Chain	Res	Type	RSRZ
2	CB	70	PHE	3.9
37	DC	177	LYS	3.9
14	CN	5	ALA	3.9
1	CA	76	C	3.9
1	CA	992	U	3.9
17	AQ	65	ILE	3.9
35	BA	2109	U	3.9
36	DB	34	U	3.9
35	DA	1526	G	3.9
37	BC	55	ASP	3.9
3	AC	64	VAL	3.9
1	CA	925	G	3.9
23	CW	60	A	3.9
36	DB	56	G	3.9
32	D6	49	HIS	3.9
35	DA	272(I)	U	3.9
40	DF	24	LEU	3.9
7	CG	76	ARG	3.9
35	BA	2190	G	3.9
35	BA	2891	G	3.9
25	AY	68	VAL	3.9
35	DA	271(J)	C	3.9
35	DA	1053	C	3.9
37	BC	45	ALA	3.9
37	DC	145	VAL	3.9
3	AC	103	VAL	3.9
38	DD	26	LYS	3.9
1	AA	1028	C	3.9
35	DA	311	A	3.9
35	DA	2801(A)	A	3.9
56	DZ	172	ALA	3.9
37	BC	99	ILE	3.9
9	CI	53	VAL	3.9
37	BC	76	ALA	3.9
26	D0	16	SER	3.9
37	BC	92	ASP	3.9
37	DC	66	HIS	3.8
35	BA	333	G	3.8
35	BA	1052	C	3.8
1	CA	1257	U	3.8
1	AA	1257	U	3.8
35	BA	2479	G	3.8

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Mol	Chain	Res	Type	RSRZ
35	BA	2794	C	3.8
35	DA	12	U	3.8
13	CM	104	ARG	3.8
39	BE	42	ASP	3.8
35	BA	476	G	3.8
36	DB	27	C	3.8
10	CJ	82	ILE	3.8
35	DA	1509(A)	A	3.8
43	BI	61	ARG	3.8
23	CW	69	C	3.8
1	AA	1047	G	3.8
20	AT	70	SER	3.8
37	DC	217	THR	3.8
23	CW	28	U	3.8
40	DF	134	GLY	3.8
35	DA	1484	G	3.8
35	DA	1586	A	3.8
2	AB	229	VAL	3.8
1	CA	1038	C	3.8
2	AB	207	ALA	3.8
41	BG	90	LEU	3.8
43	DI	101	LEU	3.8
20	CT	105	SER	3.8
1	CA	1274	G	3.8
35	BA	344	G	3.8
1	AA	81	U	3.8
25	CY	75	ALA	3.8
21	CU	24	ARG	3.8
35	BA	1214	A	3.8
1	AA	924	C	3.8
35	BA	2174	C	3.8
1	AA	1124	G	3.7
43	DI	65	ALA	3.7
25	AY	58	VAL	3.7
35	DA	2411	A	3.7
35	BA	1051	G	3.7
42	BH	137	ASP	3.7
1	CA	1044	A	3.7
37	DC	209	LEU	3.7
35	DA	506	G	3.7
39	BE	34	VAL	3.7
1	CA	921	U	3.7

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Mol	Chain	Res	Type	RSRZ
1	CA	1303	C	3.7
35	DA	1532	C	3.7
35	DA	2163	C	3.7
39	BE	35	GLN	3.7
35	BA	2123	G	3.7
35	DA	898	C	3.7
39	BE	48	GLN	3.7
1	AA	639	G	3.7
35	DA	1888	G	3.7
55	DY	54	LYS	3.7
1	CA	840	C	3.7
35	BA	1530	C	3.7
35	BA	1895	C	3.7
49	DS	56	LEU	3.7
11	CK	129	SER	3.7
23	AW	62	C	3.7
35	BA	1445(A)	C	3.7
35	BA	1863	G	3.7
35	DA	363(F)	A	3.7
42	DH	101	ARG	3.7
18	CR	87	ARG	3.7
32	D6	44	ARG	3.7
1	CA	1086	U	3.7
52	BV	5	VAL	3.7
13	AM	120	LYS	3.7
1	CA	1273	G	3.6
2	AB	222	ILE	3.6
37	DC	146	GLY	3.6
35	DA	1461	G	3.6
37	BC	192	PHE	3.6
35	BA	271(O)	C	3.6
35	BA	2172	U	3.6
1	AA	1036	G	3.6
2	CB	208	ILE	3.6
44	BN	24	GLY	3.6
47	BQ	30	GLY	3.6
1	AA	1128	C	3.6
1	AA	1270	C	3.6
7	CG	3	ARG	3.6
35	BA	1880	C	3.6
42	BH	51	ARG	3.6
37	BC	43	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
52	DV	65	GLY	3.6
1	CA	91	C	3.6
1	CA	994	A	3.6
35	BA	2128	C	3.6
1	AA	1272	G	3.6
1	CA	146	G	3.6
35	DA	271(M)	G	3.6
37	DC	184	LYS	3.6
35	DA	1113	U	3.6
23	CW	13	C	3.6
11	CK	42	TRP	3.6
35	BA	2101	G	3.6
35	DA	1170	G	3.6
35	DA	2474	C	3.6
11	CK	13	GLN	3.6
35	BA	1519	G	3.6
35	DA	407	G	3.6
1	AA	748	C	3.6
5	CE	29	GLY	3.6
35	DA	429	A	3.6
55	DY	51	VAL	3.6
21	CU	25	LYS	3.6
21	CU	23	PRO	3.6
35	DA	1173	G	3.6
1	AA	201	C	3.6
23	CW	70	C	3.5
35	BA	1881	C	3.5
9	CI	20	ARG	3.5
1	AA	72	C	3.5
1	AA	88	A	3.5
13	CM	108	ARG	3.5
35	BA	1047	G	3.5
35	BA	2127	G	3.5
35	BA	2466	C	3.5
35	DA	299	A	3.5
30	D4	23	GLU	3.5
52	BV	17	GLY	3.5
35	BA	403	U	3.5
1	AA	1314	C	3.5
35	DA	415	A	3.5
35	DA	271(L)	U	3.5
35	DA	2185	C	3.5

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Mol	Chain	Res	Type	RSRZ
37	DC	149	ILE	3.5
2	CB	215	LEU	3.5
28	D2	62	THR	3.5
1	CA	488	C	3.5
35	BA	1018	C	3.5
37	BC	103	ILE	3.5
11	AK	19	ALA	3.5
35	BA	2482	G	3.5
37	DC	92	ASP	3.5
52	BV	28	GLU	3.5
1	CA	181	G	3.5
1	CA	413	G	3.5
35	DA	495	G	3.5
35	DA	334	C	3.5
1	AA	925	G	3.5
1	CA	837	G	3.5
35	BA	1418	G	3.5
43	DI	73	GLU	3.5
17	AQ	19	VAL	3.5
1	AA	1447	A	3.5
37	BC	174	PRO	3.5
25	CY	55	ILE	3.5
25	AY	107	THR	3.5
46	DP	118	GLY	3.5
3	CC	104	GLN	3.4
43	BI	136	VAL	3.4
1	CA	1030(D)	A	3.4
19	CS	35	SER	3.4
35	DA	333	G	3.4
35	DA	1557	C	3.4
55	BY	46	LYS	3.4
7	CG	91	VAL	3.4
42	BH	52	VAL	3.4
20	CT	106	ALA	3.4
55	BY	51	VAL	3.4
1	CA	1010	G	3.4
1	CA	1140	C	3.4
35	BA	2103	C	3.4
1	AA	365	U	3.4
37	DC	107	TRP	3.4
43	DI	88	ILE	3.4
2	AB	37	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
1	CA	1395	C	3.4
22	AV	27	G	3.4
35	DA	1740	G	3.4
41	DG	74	LYS	3.4
43	DI	56	LYS	3.4
35	BA	1528	A	3.4
35	DA	900	A	3.4
36	BB	112	U	3.4
43	DI	146	ALA	3.4
49	DS	50	SER	3.4
32	B6	47	THR	3.4
1	AA	1001(A)	G	3.4
35	BA	647	G	3.4
1	AA	366	C	3.4
52	BV	46	VAL	3.4
35	BA	1494	A	3.4
35	DA	1749	A	3.4
25	AY	110	ARG	3.4
40	BF	19	GLU	3.4
46	BP	138	LEU	3.3
49	DS	48	LEU	3.3
35	DA	406	G	3.3
28	D2	31	GLU	3.3
43	DI	142	VAL	3.3
35	BA	497	A	3.3
35	BA	1106	A	3.3
35	BA	645	C	3.3
35	BA	1866	C	3.3
19	CS	68	GLY	3.3
35	BA	1954	G	3.3
36	DB	107	G	3.3
37	DC	86	ALA	3.3
10	CJ	72	VAL	3.3
35	DA	300	A	3.3
35	DA	918	A	3.3
23	CW	3	C	3.3
36	DB	3	C	3.3
37	BC	46	LYS	3.3
23	CW	71	G	3.3
35	DA	1112	G	3.3
35	BA	2171	A	3.3
1	CA	1162	C	3.3

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Mol	Chain	Res	Type	RSRZ
35	BA	2107	C	3.3
43	DI	25	TYR	3.3
16	CP	83	GLU	3.3
1	CA	183	G	3.3
37	DC	22	ILE	3.3
35	DA	1052	C	3.3
49	BS	33	LYS	3.3
9	CI	50	LEU	3.3
1	AA	160	A	3.3
1	CA	1272	G	3.3
35	BA	1042	G	3.3
35	DA	171	G	3.3
35	DA	1529	G	3.3
36	DB	110	G	3.3
43	DI	89	TYR	3.3
1	AA	156	G	3.3
35	DA	855	G	3.3
35	DA	919	G	3.3
35	DA	2313	C	3.3
35	DA	2788	C	3.3
37	DC	153	ILE	3.3
2	AB	237	ALA	3.3
35	DA	1496	A	3.3
23	CW	65	G	3.3
35	BA	1114	G	3.3
35	DA	1546	C	3.3
7	CG	90	GLU	3.3
49	DS	59	LYS	3.3
1	CA	1004	A	3.3
35	BA	2832	U	3.3
35	DA	271(N)	U	3.3
43	DI	134	PRO	3.3
35	BA	275	G	3.3
41	BG	44	GLY	3.3
37	BC	19	VAL	3.3
1	CA	1302	U	3.2
41	BG	41	GLN	3.2
35	BA	1450(A)	C	3.2
19	CS	81	ARG	3.2
20	CT	103	GLY	3.2
23	CW	19	G	3.2
25	AY	69	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
35	DA	2351	G	3.2
39	BE	204	ALA	3.2
42	DH	83	TYR	3.2
44	BN	134	ARG	3.2
1	CA	196	A	3.2
10	AJ	83	GLU	3.2
46	BP	130	PHE	3.2
26	D0	5	LYS	3.2
28	B2	35	LEU	3.2
1	CA	1229	A	3.2
35	BA	413	C	3.2
35	BA	2481	G	3.2
35	DA	356	G	3.2
37	DC	124	GLY	3.2
35	DA	1106	A	3.2
37	DC	24	GLU	3.2
37	DC	127	LEU	3.2
16	CP	13	HIS	3.2
1	CA	718	G	3.2
37	DC	76	ALA	3.2
1	AA	1004	A	3.2
47	BQ	138	ASP	3.2
23	CW	29	C	3.2
35	BA	2163	C	3.2
3	CC	80	GLY	3.2
25	CY	99	LEU	3.2
35	BA	602	G	3.2
35	DA	2104	G	3.2
36	DB	15	A	3.2
13	AM	6	GLY	3.2
42	BH	34	GLU	3.2
46	DP	84	ASN	3.2
1	AA	700	G	3.2
35	BA	351	G	3.2
35	BA	707	G	3.2
35	DA	881	G	3.2
35	DA	1107	G	3.2
35	DA	1421	G	3.2
5	CE	113	ALA	3.2
13	AM	19	LEU	3.2
42	BH	58	GLU	3.2
7	CG	77	SER	3.2

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Mol	Chain	Res	Type	RSRZ
13	CM	125	ARG	3.2
35	BA	1544	A	3.2
56	DZ	105	VAL	3.2
43	DI	80	PRO	3.2
35	BA	2344	U	3.1
41	DG	58	GLN	3.1
37	DC	185	LEU	3.1
35	DA	1416	G	3.1
1	AA	1223	C	3.1
35	BA	1145	C	3.1
42	BH	138	LYS	3.1
19	AS	81	ARG	3.1
50	DT	115	ARG	3.1
41	DG	32	PRO	3.1
1	AA	1446	U	3.1
43	BI	131	LYS	3.1
17	CQ	18	THR	3.1
25	AY	57	THR	3.1
35	DA	289	A	3.1
1	AA	272	C	3.1
36	DB	24	G	3.1
7	AG	49	ILE	3.1
35	DA	2172	U	3.1
37	DC	77	ILE	3.1
1	CA	177	C	3.1
35	DA	1506	C	3.1
42	BH	139	GLN	3.1
35	DA	1046	A	3.1
7	CG	89	MET	3.1
35	DA	634	C	3.1
25	AY	98	ALA	3.1
35	BA	2807	G	3.1
35	DA	1507	A	3.1
37	BC	75	LEU	3.1
43	DI	116	LEU	3.1
43	BI	72	LEU	3.1
1	CA	78	G	3.1
35	BA	2661	G	3.1
35	DA	1176	G	3.1
35	DA	1533	G	3.1
41	BG	92	VAL	3.1
1	AA	430	A	3.1

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Mol	Chain	Res	Type	RSRZ
1	CA	1236	A	3.1
35	BA	896	A	3.1
35	BA	2809	A	3.1
35	DA	1587	A	3.1
36	DB	53	A	3.1
37	BC	156	ILE	3.1
35	BA	1463	C	3.1
42	BH	50	VAL	3.1
5	CE	83	GLU	3.1
1	CA	1304	G	3.1
35	DA	363(B)	G	3.1
35	DA	600	G	3.1
35	DA	934	G	3.1
25	AY	70	SER	3.1
7	AG	154	TYR	3.1
29	B3	1	MET	3.1
1	CA	1224	G	3.1
25	CY	41	LEU	3.1
35	BA	1950	G	3.1
44	BN	127	ASP	3.1
17	CQ	45	HIS	3.1
35	DA	1499	C	3.1
2	AB	99	GLY	3.1
1	AA	1024	G	3.1
1	AA	1068	G	3.1
23	AW	10	G	3.1
35	DA	2112	G	3.1
8	AH	35	ILE	3.1
1	AA	421	U	3.0
3	AC	65	ALA	3.0
35	DA	2312	U	3.0
42	BH	66	GLY	3.0
34	D8	33	ASN	3.0
37	DC	207	THR	3.0
37	DC	50	ASP	3.0
1	CA	414	A	3.0
52	BV	96	ILE	3.0
30	D4	3	GLU	3.0
37	DC	98	GLU	3.0
37	DC	171	ILE	3.0
43	DI	78	THR	3.0
52	DV	96	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	AA	991	U	3.0
1	CA	1125	U	3.0
35	DA	1528(A)	A	3.0
35	DA	2327	A	3.0
38	BD	93	ALA	3.0
49	DS	37	ALA	3.0
2	AB	76	GLN	3.0
1	AA	144	G	3.0
35	BA	271(M)	G	3.0
35	BA	882	G	3.0
35	DA	2100	G	3.0
35	BA	280	C	3.0
39	DE	69	LYS	3.0
8	CH	130	GLY	3.0
1	CA	1121	U	3.0
6	AF	39	LYS	3.0
25	AY	152	ASP	3.0
1	CA	1022	G	3.0
35	BA	2526	G	3.0
35	DA	2811	G	3.0
37	DC	87	GLU	3.0
43	BI	10	GLU	3.0
23	CW	44	A	3.0
30	D4	9	LEU	3.0
1	AA	418	C	3.0
1	CA	1320	C	3.0
20	CT	89	ARG	3.0
56	BZ	89	PHE	3.0
25	CY	69	GLN	3.0
1	AA	1126	U	3.0
25	CY	144	ALA	3.0
33	D7	49	ARG	3.0
10	CJ	35	SER	3.0
17	CQ	65	ILE	3.0
1	AA	200	G	3.0
1	AA	1037	C	3.0
10	CJ	70	ARG	3.0
17	AQ	44	ALA	3.0
35	DA	176	G	3.0
35	DA	883	G	3.0
9	CI	65	VAL	3.0
35	DA	913	U	3.0

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Mol	Chain	Res	Type	RSRZ
14	AN	2	ALA	3.0
52	BV	30	GLY	3.0
9	CI	14	VAL	3.0
35	BA	34	C	3.0
35	BA	1509	C	3.0
35	DA	1169	G	3.0
32	D6	41	PRO	3.0
35	BA	363(F)	A	3.0
50	BT	136	GLN	3.0
32	D6	13	CYS	3.0
1	AA	1061	G	3.0
1	CA	428	G	3.0
35	BA	1448	G	3.0
35	BA	1921	G	3.0
35	DA	360	G	3.0
52	DV	19	LYS	3.0
11	CK	44	SER	3.0
55	BY	59	GLY	3.0
26	D0	17	GLN	3.0
37	BC	93	TYR	3.0
1	CA	268	C	2.9
41	DG	73	ALA	2.9
49	DS	104	GLY	2.9
1	CA	77	G	2.9
1	CA	1042	G	2.9
35	BA	1465	G	2.9
37	DC	126	LYS	2.9
43	BI	68	LEU	2.9
1	AA	994	A	2.9
13	AM	64	TRP	2.9
37	BC	71	GLN	2.9
41	BG	134	GLY	2.9
37	DC	189	ILE	2.9
35	BA	1108	U	2.9
35	BA	894	C	2.9
35	BA	1109	C	2.9
9	CI	94	ALA	2.9
37	BC	89	ALA	2.9
1	AA	69	G	2.9
1	AA	1130	A	2.9
35	BA	1001	A	2.9
35	BA	1533	G	2.9

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Mol	Chain	Res	Type	RSRZ
35	DA	2472	G	2.9
16	AP	84	ALA	2.9
1	AA	1001	A	2.9
1	CA	530	G	2.9
35	DA	329	G	2.9
35	DA	878	A	2.9
35	DA	1303	G	2.9
56	DZ	104	PHE	2.9
13	AM	122	LYS	2.9
35	BA	1166	C	2.9
35	DA	2128	C	2.9
46	BP	90	ARG	2.9
51	DU	77	SER	2.9
1	CA	692	U	2.9
35	DA	2119	A	2.9
35	BA	919	G	2.9
35	BA	1702	G	2.9
35	DA	175	G	2.9
35	DA	430	G	2.9
35	DA	2470	G	2.9
1	AA	1007	C	2.9
35	BA	402	A	2.9
35	DA	1528	A	2.9
42	BH	49	VAL	2.9
37	DC	148	ASN	2.9
43	BI	132	PRO	2.9
35	DA	2187	G	2.9
46	BP	118	GLY	2.9
1	AA	1172	C	2.9
1	CA	1264	C	2.9
35	BA	2551	C	2.9
35	DA	2103	C	2.9
39	BE	69	LYS	2.9
54	BX	69	TYR	2.9
23	CW	73	A	2.9
49	DS	11	LYS	2.9
1	AA	254	G	2.9
1	CA	199	G	2.9
1	CA	266	G	2.9
2	CB	120	ALA	2.9
3	CC	79	ARG	2.9
35	BA	628	G	2.9

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Mol	Chain	Res	Type	RSRZ
1	AA	91	C	2.9
52	BV	98	GLU	2.9
1	AA	1020	U	2.9
35	BA	1639	U	2.9
35	BA	1940	U	2.9
25	AY	45	TYR	2.9
19	CS	72	GLY	2.9
40	BF	11	VAL	2.9
49	DS	38	GLN	2.9
1	CA	157	G	2.9
1	CA	1178	G	2.9
35	BA	1169	G	2.9
35	BA	1667	G	2.9
44	BN	26	LEU	2.9
35	BA	1113	U	2.9
35	DA	2179	C	2.9
35	BA	918	A	2.9
1	CA	995	C	2.9
35	BA	1030	G	2.9
35	BA	2472	G	2.9
35	DA	414	C	2.9
35	DA	1042	G	2.9
35	DA	2796	U	2.9
41	DG	80	PHE	2.9
42	BH	47	GLU	2.8
41	DG	81	LYS	2.8
23	AW	21	U	2.8
49	BS	49	VAL	2.8
1	CA	254	G	2.8
35	BA	1638	C	2.8
10	AJ	35	SER	2.8
37	DC	105	ASP	2.8
56	BZ	7	ALA	2.8
35	DA	1213	A	2.8
35	DA	1876	A	2.8
17	CQ	63	ARG	2.8
1	CA	571	U	2.8
49	DS	31	SER	2.8
3	CC	36	ASP	2.8
17	CQ	70	ARG	2.8
23	AW	29	C	2.8
35	BA	1498	C	2.8

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Mol	Chain	Res	Type	RSRZ
55	DY	61	ILE	2.8
7	CG	75	VAL	2.8
37	BC	59	ARG	2.8
40	BF	27	GLU	2.8
9	CI	36	TYR	2.8
35	BA	2538	C	2.8
35	DA	271(O)	C	2.8
35	DA	1967	C	2.8
37	BC	26	ALA	2.8
26	B0	16	SER	2.8
1	AA	98	G	2.8
35	BA	1888	G	2.8
35	DA	140	G	2.8
35	DA	1579	A	2.8
36	DB	26	A	2.8
3	AC	100	ALA	2.8
25	AY	60	ALA	2.8
52	DV	45	THR	2.8
41	BG	78	SER	2.8
9	AI	19	LEU	2.8
25	AY	66	LEU	2.8
37	BC	97	GLU	2.8
37	BC	153	ILE	2.8
1	AA	269	C	2.8
35	BA	343	C	2.8
1	CA	953	G	2.8
17	AQ	43	LEU	2.8
35	BA	629	G	2.8
49	DS	52	SER	2.8
9	CI	60	ASP	2.8
3	CC	78	GLY	2.8
1	AA	1264	C	2.8
42	BH	35	VAL	2.8
35	BA	1579	A	2.8
35	DA	2883	A	2.8
51	BU	91	ASP	2.8
55	DY	28	LYS	2.8
1	CA	1370	G	2.8
35	BA	1435	G	2.8
41	DG	34	LEU	2.8
7	AG	57	GLU	2.8
35	DA	184	C	2.8

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Mol	Chain	Res	Type	RSRZ
51	DU	74	LEU	2.8
37	DC	188	ASN	2.8
17	CQ	99	SER	2.8
37	BC	221	SER	2.8
37	DC	44	HIS	2.8
1	CA	286	G	2.8
7	AG	62	PHE	2.8
37	BC	22	ILE	2.8
1	AA	417	C	2.8
1	AA	419	C	2.8
25	AY	185	GLY	2.8
35	BA	2480	C	2.8
43	BI	65	ALA	2.8
2	CB	68	ILE	2.8
9	CI	66	ARG	2.8
23	CW	43	G	2.8
37	DC	83	ILE	2.8
35	BA	288	C	2.8
35	DA	1638	C	2.8
9	CI	18	PHE	2.8
35	BA	1509(B)	A	2.8
35	BA	1876	A	2.8
1	CA	156	G	2.7
1	CA	265	G	2.7
9	CI	3	GLN	2.7
25	AY	56	ALA	2.7
1	CA	1037	C	2.7
35	BA	1147	C	2.7
49	DS	34	HIS	2.7
35	BA	501	A	2.7
35	BA	1637	A	2.7
35	BA	1212	G	2.7
35	BA	1552	G	2.7
35	DA	1108	U	2.7
1	CA	419	C	2.7
35	BA	1181	C	2.7
17	CQ	17	LYS	2.7
35	DA	332	A	2.7
44	BN	23	LEU	2.7
13	AM	24	GLY	2.7
37	BC	158	ALA	2.7
39	BE	205	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
40	BF	23	ASP	2.7
46	BP	119	GLU	2.7
1	AA	1032	G	2.7
35	BA	1532	C	2.7
35	BA	2801(A)	A	2.7
1	CA	96	U	2.7
39	DE	90	THR	2.7
11	CK	17	GLY	2.7
1	AA	68	G	2.7
35	BA	407	G	2.7
35	BA	2805	G	2.7
36	BB	110	G	2.7
1	AA	431	A	2.7
7	AG	4	ARG	2.7
20	AT	98	PRO	2.7
23	AW	36	A	2.7
4	AD	25	ARG	2.7
11	AK	43	SER	2.7
35	DA	1402	C	2.7
35	DA	1837	C	2.7
3	CC	90	GLU	2.7
13	CM	61	GLU	2.7
1	CA	189(B)	C	2.7
41	DG	122	PRO	2.7
46	BP	114	ILE	2.7
1	CA	1225	A	2.7
3	AC	35	GLU	2.7
35	BA	175	G	2.7
35	DA	1418	G	2.7
13	CM	3	ARG	2.7
28	D2	34	GLU	2.7
1	AA	1161	C	2.7
3	CC	98	ASN	2.7
1	CA	1095	U	2.7
35	BA	1213	A	2.7
4	CD	40	PRO	2.7
10	CJ	85	LEU	2.7
35	BA	271(R)	G	2.7
35	BA	1922	G	2.7
35	DA	602	G	2.7
43	DI	105	HIS	2.7
35	DA	1509	C	2.7

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Mol	Chain	Res	Type	RSRZ
35	DA	1497	U	2.7
41	BG	159	VAL	2.7
2	CB	42	ILE	2.7
1	AA	460	G	2.7
1	CA	1036	G	2.7
37	DC	142	ALA	2.7
9	CI	2	GLU	2.6
25	CY	68	VAL	2.6
25	AY	54	GLN	2.6
56	DZ	179	ASP	2.6
26	B0	8	ALA	2.6
55	BY	47	LYS	2.6
35	DA	1866	C	2.6
35	DA	1895	C	2.6
36	BB	27	C	2.6
3	CC	68	VAL	2.6
7	AG	68	ASN	2.6
36	DB	48	A	2.6
23	AW	31	G	2.6
35	BA	308	G	2.6
35	BA	1630	G	2.6
25	AY	17	SER	2.6
25	CY	52	LEU	2.6
35	DA	603	A	2.6
37	DC	175	VAL	2.6
35	BA	1154	G	2.6
35	BA	1478	G	2.6
1	AA	1430	C	2.6
1	CA	188	C	2.6
50	DT	119	LYS	2.6
17	AQ	101	ARG	2.6
35	BA	345	A	2.6
35	BA	429	A	2.6
35	DA	2654	A	2.6
13	AM	54	VAL	2.6
41	BG	182	LYS	2.6
41	DG	35	GLU	2.6
1	CA	1023	G	2.6
35	BA	1878	G	2.6
23	CW	16	C	2.6
35	DA	1710	C	2.6
37	BC	212	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
35	BA	502	A	2.6
35	BA	1936	A	2.6
35	DA	2629	A	2.6
42	BH	45	VAL	2.6
1	CA	64	G	2.6
23	AW	5	G	2.6
23	CW	31	G	2.6
35	BA	75	G	2.6
36	BB	63	G	2.6
56	BZ	4	ARG	2.6
13	CM	105	THR	2.6
25	CY	66	LEU	2.6
3	CC	99	VAL	2.6
25	CY	61	PRO	2.6
35	DA	2188	C	2.6
35	DA	2896	C	2.6
35	BA	1545	A	2.6
9	CI	6	GLY	2.6
35	DA	1341	U	2.6
35	BA	15	G	2.6
35	BA	2831	G	2.6
1	CA	47	C	2.6
1	CA	924	C	2.6
46	DP	87	ASP	2.6
1	AA	751	U	2.6
41	BG	156	ASP	2.6
1	AA	1266	G	2.6
1	AA	1322	C	2.6
1	CA	1174	G	2.6
10	CJ	97	GLU	2.6
35	BA	892	G	2.6
35	BA	1144	G	2.6
35	BA	1447	G	2.6
35	DA	258	G	2.6
35	BA	2533	A	2.6
43	DI	77	LEU	2.6
1	CA	1122	U	2.6
56	BZ	104	PHE	2.6
37	DC	84	LYS	2.6
49	DS	44	LYS	2.6
1	AA	179	A	2.6
1	AA	1504	G	2.6

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Mol	Chain	Res	Type	RSRZ
35	DA	272(B)	G	2.6
56	DZ	139	VAL	2.5
1	CA	1046	A	2.5
2	AB	38	GLY	2.5
3	CC	74	GLY	2.5
11	CK	46	GLY	2.5
26	B0	85	ALA	2.5
35	BA	2180	U	2.5
36	BB	79	C	2.5
35	DA	2810	A	2.5
37	DC	73	ARG	2.5
1	AA	1127	G	2.5
1	AA	1474	G	2.5
35	DA	2782	G	2.5
36	DB	61	G	2.5
37	BC	42	GLU	2.5
56	BZ	12	GLY	2.5
20	CT	92	LEU	2.5
52	BV	97	LYS	2.5
1	AA	217	C	2.5
1	AA	429	U	2.5
35	BA	1178	C	2.5
9	AI	111	ARG	2.5
9	CI	34	ASN	2.5
15	AO	22	THR	2.5
35	DA	629	G	2.5
35	DA	1047	G	2.5
35	DA	2894	G	2.5
25	AY	67	VAL	2.5
35	BA	2748	A	2.5
36	DB	66	A	2.5
41	DG	78	SER	2.5
1	AA	1461	G	2.5
25	AY	61	PRO	2.5
26	B0	17	GLN	2.5
32	D6	45	LYS	2.5
35	DA	1388	G	2.5
35	DA	1949	G	2.5
35	DA	2659	G	2.5
45	BO	59	LYS	2.5
32	B6	26	ASN	2.5
52	DV	20	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	AA	1219	U	2.5
37	DC	155	GLU	2.5
25	AY	48	ALA	2.5
37	BC	219	GLY	2.5
47	BQ	124	LYS	2.5
35	BA	256	A	2.5
35	DA	1877	A	2.5
3	AC	62	ASP	2.5
13	AM	108	ARG	2.5
1	AA	1265	G	2.5
1	CA	1082	G	2.5
11	AK	16	SER	2.5
35	BA	1107	G	2.5
51	BU	118	GLY	2.5
13	CM	25	ILE	2.5
35	BA	2790	A	2.5
56	DZ	59	LEU	2.5
7	AG	75	VAL	2.5
17	AQ	68	ARG	2.5
1	AA	384	G	2.5
1	CA	1392	G	2.5
23	AW	48	U	2.5
4	CD	44	GLY	2.5
35	BA	217	G	2.5
35	BA	2782	G	2.5
1	CA	1116	C	2.5
1	CA	1244	C	2.5
39	BE	49	LEU	2.5
12	CL	33	ARG	2.5
36	BB	77	U	2.5
52	BV	29	PRO	2.5
35	BA	1416	G	2.5
35	DA	1647	G	2.5
3	CC	64	VAL	2.5
26	D0	61	ALA	2.5
1	AA	47	C	2.5
1	AA	689	C	2.5
1	CA	695	A	2.5
7	AG	5	ARG	2.5
37	BC	148	ASN	2.5
56	DZ	96	VAL	2.5
35	BA	1557	C	2.5

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Mol	Chain	Res	Type	RSRZ
35	DA	257	A	2.5
35	DA	605	C	2.5
35	DA	1467	C	2.5
49	DS	54	LEU	2.5
43	DI	74	ASN	2.5
1	CA	73	G	2.5
10	CJ	91	PRO	2.5
35	BA	2468	G	2.5
35	BA	2484	G	2.5
35	DA	2833	G	2.5
1	AA	271	C	2.5
1	AA	385	C	2.5
1	CA	983	A	2.5
17	CQ	68	ARG	2.5
35	BA	105	C	2.5
40	DF	1	MET	2.5
35	BA	406	G	2.4
35	BA	2861	G	2.4
35	DA	864	G	2.4
35	DA	1904	G	2.4
7	CG	13	GLN	2.4
7	CG	16	LEU	2.4
11	AK	31	THR	2.4
35	DA	1109	C	2.4
18	AR	61	LYS	2.4
35	BA	895	U	2.4
37	BC	157	LYS	2.4
50	DT	1	MET	2.4
52	DV	31	ALA	2.4
1	AA	589	C	2.4
1	AA	1300	G	2.4
1	CA	144	G	2.4
1	CA	1139	G	2.4
35	BA	875	G	2.4
35	BA	1470	G	2.4
35	BA	2757	A	2.4
35	DA	361	G	2.4
35	DA	1049	C	2.4
35	DA	1648	C	2.4
35	DA	2805	G	2.4
35	DA	2807	G	2.4
36	DB	64	C	2.4

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Mol	Chain	Res	Type	RSRZ
52	BV	43	GLU	2.4
2	AB	91	PRO	2.4
2	AB	209	ARG	2.4
2	CB	76	GLN	2.4
26	D0	15	ASP	2.4
28	D2	41	ILE	2.4
37	BC	139	ASN	2.4
37	DC	136	LEU	2.4
37	DC	212	VAL	2.4
41	BG	91	ARG	2.4
1	AA	601	C	2.4
1	AA	1370	G	2.4
1	CA	143	A	2.4
1	CA	184	G	2.4
1	CA	1032	G	2.4
1	CA	1117	G	2.4
1	CA	1138	G	2.4
35	BA	1116	C	2.4
35	BA	1586	A	2.4
35	DA	268	C	2.4
35	DA	271(K)	U	2.4
37	BC	27	ARG	2.4
41	BG	81	LYS	2.4
41	DG	50	ALA	2.4
37	DC	54	SER	2.4
32	D6	12	GLU	2.4
1	CA	562	C	2.4
23	AW	17	C	2.4
35	BA	603	A	2.4
35	BA	2117	A	2.4
35	DA	924	C	2.4
1	CA	947	G	2.4
35	DA	707	G	2.4
35	DA	1212	G	2.4
35	DA	1482	G	2.4
35	DA	2165	G	2.4
37	BC	191	ALA	2.4
43	BI	54	GLN	2.4
19	CS	18	LYS	2.4
19	CS	36	ARG	2.4
32	D6	14	THR	2.4
52	BV	26	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
44	DN	13	TRP	2.4
55	DY	83	THR	2.4
42	BH	33	LEU	2.4
1	AA	153	C	2.4
1	CA	1218	C	2.4
35	BA	1464	C	2.4
35	BA	2474	C	2.4
35	DA	477	A	2.4
35	DA	2863	C	2.4
37	DC	125	SER	2.4
45	BO	57	VAL	2.4
1	AA	145	G	2.4
1	CA	1011	G	2.4
1	CA	1030(A)	G	2.4
35	BA	354	G	2.4
35	BA	881	G	2.4
35	BA	1421	G	2.4
35	DA	892	G	2.4
7	AG	53	LYS	2.4
32	D6	37	ARG	2.4
41	BG	155	MET	2.4
50	BT	2	ASN	2.4
18	CR	85	LEU	2.4
35	DA	1498	C	2.4
1	AA	1190	G	2.4
13	AM	116	THR	2.4
31	B5	54	GLY	2.4
35	BA	317	G	2.4
35	BA	458	G	2.4
37	DC	79	LYS	2.4
1	AA	480	U	2.4
1	AA	1052	U	2.4
25	CY	76	LEU	2.4
35	DA	1523	U	2.4
2	AB	32	ILE	2.4
35	DA	319	C	2.4
35	DA	1320	C	2.4
37	DC	173	ALA	2.4
46	DP	96	THR	2.4
37	BC	145	VAL	2.4
1	AA	117	G	2.4
1	AA	424	G	2.4

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Mol	Chain	Res	Type	RSRZ
1	CA	301	G	2.4
11	AK	29	ILE	2.4
35	DA	158	U	2.4
43	BI	138	ILE	2.4
7	CG	73	MET	2.4
37	DC	80	GLY	2.4
5	AE	120	THR	2.4
13	AM	50	GLU	2.4
35	BA	1631	C	2.4
35	BA	2568	C	2.4
35	DA	646	A	2.4
35	DA	1469	A	2.4
35	DA	2666	C	2.4
52	BV	99	ILE	2.4
7	AG	2	ALA	2.4
1	CA	592	G	2.4
20	CT	98	PRO	2.4
35	BA	1466	G	2.4
2	AB	85	ALA	2.4
49	BS	48	LEU	2.4
49	DS	27	SER	2.4
1	AA	1158	C	2.4
1	AA	1321	C	2.4
35	BA	817	C	2.4
35	DA	416	C	2.4
2	CB	96	ARG	2.3
35	BA	157	U	2.3
11	AK	82	VAL	2.3
11	CK	82	VAL	2.3
43	BI	126	TYR	2.3
49	BS	11	LYS	2.3
1	CA	1177	G	2.3
35	BA	430	G	2.3
35	DA	1642	G	2.3
25	CY	71	TRP	2.3
39	DE	73	GLU	2.3
52	BV	100	ARG	2.3
1	AA	1363(A)	A	2.3
1	CA	1447	A	2.3
35	DA	1027	A	2.3
42	BH	20	ALA	2.3
1	AA	1431	C	2.3

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Mol	Chain	Res	Type	RSRZ
1	CA	591	U	2.3
35	BA	1911	U	2.3
7	AG	155	ARG	2.3
51	BU	84	LYS	2.3
1	CA	691	G	2.3
35	BA	1529	G	2.3
35	BA	2663	G	2.3
1	AA	704	A	2.3
32	B6	35	GLU	2.3
14	AN	61	TRP	2.3
35	BA	257	A	2.3
35	BA	2411	A	2.3
35	BA	2835	A	2.3
35	DA	1449	A	2.3
35	DA	1509(B)	A	2.3
42	DH	34	GLU	2.3
1	AA	1393	U	2.3
1	CA	1226	C	2.3
37	DC	68	LEU	2.3
46	DP	142	GLY	2.3
49	BS	32	LEU	2.3
9	CI	21	PRO	2.3
17	CQ	43	LEU	2.3
41	DG	43	LEU	2.3
44	DN	139	GLU	2.3
25	AY	75	ALA	2.3
46	BP	113	LYS	2.3
1	CA	923	A	2.3
1	CA	1084	G	2.3
9	CI	93	ARG	2.3
36	DB	23	G	2.3
34	B8	31	HIS	2.3
44	BN	30	ILE	2.3
1	AA	600	C	2.3
1	AA	999	C	2.3
1	CA	1029	C	2.3
19	CS	71	LEU	2.3
35	BA	1582	C	2.3
41	DG	172	LEU	2.3
49	DS	51	ALA	2.3
3	CC	207	VAL	2.3
11	AK	26	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	CA	603	U	2.3
1	AA	155	C	2.3
1	AA	266	G	2.3
1	AA	840	C	2.3
1	AA	1181	G	2.3
35	DA	523	C	2.3
35	DA	635	C	2.3
35	DA	1883	G	2.3
36	DB	5	C	2.3
7	AG	48	LYS	2.3
13	AM	126	LYS	2.3
37	BC	18	LYS	2.3
43	DI	68	LEU	2.3
37	DC	191	ALA	2.3
35	DA	1757	U	2.3
1	CA	189(L)	G	2.3
1	CA	285	G	2.3
16	CP	84	ALA	2.3
17	CQ	101	ARG	2.3
32	B6	29	ASN	2.3
35	BA	2896	C	2.3
3	CC	91	LEU	2.3
41	BG	126	ASP	2.3
2	CB	163	PHE	2.3
4	AD	37	PRO	2.3
13	CM	69	GLU	2.3
35	BA	2392	A	2.3
35	BA	2602	A	2.3
35	DA	363(A)	A	2.3
35	DA	2171	A	2.3
36	BB	48	A	2.3
21	CU	2	GLY	2.3
35	BA	1146	C	2.3
1	AA	1271	G	2.3
23	AW	27	G	2.3
35	DA	1878	G	2.3
46	DP	120	ALA	2.3
41	DG	39	ILE	2.3
1	CA	1390	U	2.3
35	BA	2537	U	2.3
35	DA	2878	U	2.3
42	BH	74	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
50	DT	104	ASN	2.3
1	AA	224	C	2.3
11	CK	14	VAL	2.3
17	AQ	66	SER	2.3
1	AA	1082	G	2.3
1	CA	1001(A)	G	2.3
1	CA	1047	G	2.3
35	DA	1414	G	2.3
36	BB	111	G	2.3
20	AT	106	ALA	2.3
29	D3	25	ALA	2.3
1	CA	1313	U	2.3
35	BA	434	U	2.3
17	CQ	16	GLN	2.3
49	DS	103	GLU	2.3
1	AA	1225	A	2.3
19	AS	18	LYS	2.3
26	D0	4	LYS	2.3
35	BA	1153	C	2.3
35	DA	151	C	2.3
1	AA	1009	G	2.2
1	CA	198	G	2.2
1	CA	1187	G	2.2
16	AP	41	PRO	2.2
35	DA	363(C)	G	2.2
35	DA	1578	U	2.2
35	DA	2180	U	2.2
37	DC	69	GLY	2.2
41	BG	79	ASN	2.2
25	AY	59	THR	2.2
39	BE	44	TYR	2.2
35	DA	374	A	2.2
35	DA	1301	A	2.2
35	DA	1494	A	2.2
36	BB	105	A	2.2
1	CA	312	C	2.2
35	BA	2178	C	2.2
50	DT	137	LYS	2.2
37	BC	86	ALA	2.2
1	AA	1368	G	2.2
3	AC	143	GLU	2.2
56	DZ	69	THR	2.2

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Mol	Chain	Res	Type	RSRZ
11	AK	30	VAL	2.2
13	AM	104	ARG	2.2
35	DA	1286	A	2.2
1	CA	417	C	2.2
1	CA	1249	C	2.2
35	DA	2350	C	2.2
37	BC	126	LYS	2.2
1	AA	1301	U	2.2
25	AY	49	HIS	2.2
42	BH	26	VAL	2.2
1	CA	93	G	2.2
1	CA	1175	G	2.2
1	CA	1190	G	2.2
1	CA	1265	G	2.2
35	BA	1299	G	2.2
35	BA	1423	G	2.2
35	BA	1525	G	2.2
35	BA	1896	G	2.2
35	DA	259	G	2.2
35	DA	1492	G	2.2
32	D6	26	ASN	2.2
41	BG	40	ASN	2.2
42	DH	86	GLU	2.2
13	CM	66	LEU	2.2
43	BI	139	GLN	2.2
53	DW	112	GLY	2.2
35	BA	216	A	2.2
35	DA	2662	A	2.2
44	DN	14	VAL	2.2
10	AJ	70	ARG	2.2
35	DA	271(P)	C	2.2
35	DA	288	C	2.2
47	DQ	80	GLU	2.2
9	CI	7	THR	2.2
9	CI	38	GLN	2.2
1	AA	189(L)	G	2.2
1	CA	491	G	2.2
2	AB	75	LYS	2.2
49	BS	12	PHE	2.2
2	CB	33	TYR	2.2
23	AW	71	G	2.2
35	DA	281	G	2.2

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Mol	Chain	Res	Type	RSRZ
35	DA	2864	G	2.2
1	CA	313	A	2.2
17	AQ	49	GLU	2.2
35	DA	863	A	2.2
2	AB	68	ILE	2.2
35	DA	2344	U	2.2
1	AA	1230	C	2.2
7	CG	62	PHE	2.2
35	BA	1411	C	2.2
36	DB	20	C	2.2
43	DI	98	ALA	2.2
51	DU	101	ARG	2.2
11	AK	35	PRO	2.2
19	CS	37	ARG	2.2
1	AA	77	G	2.2
35	DA	88	G	2.2
55	DY	92	ASN	2.2
35	BA	1864	U	2.2
35	DA	2466	C	2.2
42	BH	56	SER	2.2
41	DG	72	ARG	2.2
51	DU	73	GLY	2.2
28	B2	38	GLN	2.2
43	DI	91	SER	2.2
56	BZ	31	ARG	2.2
1	AA	1268	A	2.2
1	AA	1433	A	2.2
35	BA	2884	U	2.2
35	BA	2802	G	2.2
35	DA	599	G	2.2
35	DA	2352	A	2.2
38	DD	107	ALA	2.2
35	BA	248	G	2.2
35	BA	880	G	2.2
36	DB	98	G	2.2
36	DB	101	G	2.2
10	CJ	74	ILE	2.2
13	CM	47	ASP	2.2
40	BF	167	ALA	2.2
46	BP	120	ALA	2.2
37	DC	131	LEU	2.2
9	CI	80	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
34	D8	28	GLY	2.2
37	BC	21	THR	2.2
1	CA	1002	G	2.2
35	BA	1717	G	2.2
35	DA	1459	G	2.2
36	DB	6	C	2.2
2	CB	119	GLU	2.2
49	DS	76	LYS	2.2
25	CY	47	GLY	2.2
3	AC	95	THR	2.2
35	DA	1639	U	2.2
46	BP	127	ALA	2.2
1	CA	130	A	2.2
35	BA	412	A	2.2
35	BA	988	A	2.2
5	CE	6	PHE	2.2
40	DF	14	PRO	2.2
1	CA	1068	G	2.2
35	BA	874	G	2.2
35	BA	2280	G	2.2
35	BA	2409	G	2.2
35	DA	1413	G	2.2
56	DZ	118	GLN	2.2
2	CB	228	GLY	2.1
1	CA	733	A	2.1
35	BA	960	A	2.1
35	DA	454	A	2.1
35	DA	643	A	2.1
35	DA	1412	A	2.1
6	AF	64	GLN	2.1
32	B6	13	CYS	2.1
1	AA	1224	G	2.1
1	CA	345	C	2.1
35	BA	1628	G	2.1
35	BA	2540	C	2.1
35	DA	2087	G	2.1
35	DA	2164	C	2.1
35	DA	2410	G	2.1
35	DA	2480	C	2.1
49	DS	68	GLN	2.1
27	B1	78	LYS	2.1
51	BU	74	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	CB	80	ILE	2.1
35	DA	633	A	2.1
27	D1	50	ARG	2.1
44	DN	15	LEU	2.1
13	AM	25	ILE	2.1
35	DA	2261	C	2.1
1	AA	90	U	2.1
1	AA	1030(C)	G	2.1
1	AA	1150	U	2.1
1	CA	473	G	2.1
35	BA	1171	G	2.1
35	BA	1173	G	2.1
35	BA	2382	G	2.1
2	AB	34	ALA	2.1
52	DV	21	ARG	2.1
9	CI	81	ILE	2.1
11	AK	42	TRP	2.1
25	CY	148	HIS	2.1
46	DP	104	GLY	2.1
32	D6	38	LYS	2.1
49	BS	62	LYS	2.1
1	AA	188	C	2.1
1	CA	1030	C	2.1
1	CA	1132	C	2.1
5	AE	69	VAL	2.1
36	DB	70	C	2.1
37	BC	154	ARG	2.1
50	BT	39	ARG	2.1
1	AA	199	G	2.1
1	AA	688	G	2.1
1	CA	1134	G	2.1
20	AT	51	GLU	2.1
35	BA	599	G	2.1
35	DA	10	G	2.1
37	BC	146	GLY	2.1
16	CP	45	THR	2.1
37	DC	102	LYS	2.1
10	AJ	90	LEU	2.1
19	CS	15	LEU	2.1
37	DC	26	ALA	2.1
37	DC	223	ARG	2.1
39	DE	68	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
2	CB	214	ILE	2.1
35	BA	1045	A	2.1
35	BA	2781	A	2.1
1	AA	427	U	2.1
1	CA	1223	C	2.1
9	CI	57	GLY	2.1
35	BA	2258	C	2.1
35	DA	198	C	2.1
16	CP	38	TYR	2.1
35	DA	1325	G	2.1
1	AA	1093	A	2.1
1	AA	1102	A	2.1
1	CA	1286	A	2.1
4	AD	174	LEU	2.1
55	DY	86	ARG	2.1
11	AK	41	THR	2.1
5	AE	70	PRO	2.1
53	DW	63	ASP	2.1
1	AA	54	C	2.1
1	CA	366	C	2.1
1	CA	1096	C	2.1
13	CM	45	VAL	2.1
37	DC	19	VAL	2.1
1	AA	428	G	2.1
1	AA	1048	G	2.1
17	CQ	20	THR	2.1
35	DA	250	G	2.1
35	DA	1525	G	2.1
35	DA	1921	G	2.1
37	DC	161	ILE	2.1
39	DE	44	TYR	2.1
1	AA	1125	U	2.1
7	CG	17	VAL	2.1
35	BA	477	A	2.1
1	CA	1008	C	2.1
35	BA	1180	C	2.1
25	AY	76	LEU	2.1
5	CE	19	MET	2.1
18	CR	50	ILE	2.1
35	DA	880	G	2.1
10	AJ	79	ARG	2.1
35	BA	2469	A	2.1

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Mol	Chain	Res	Type	RSRZ
42	DH	137	ASP	2.1
1	CA	1115	C	2.1
35	BA	366	C	2.1
35	BA	1499	C	2.1
35	BA	1584	C	2.1
35	BA	2403	C	2.1
35	DA	1464	C	2.1
35	DA	1551	C	2.1
36	BB	91	C	2.1
39	BE	68	ALA	2.1
3	CC	190	ARG	2.1
42	DH	59	ARG	2.1
5	CE	35	GLY	2.1
13	AM	8	GLU	2.1
1	CA	412	A	2.1
23	AW	28	U	2.1
35	BA	1026	U	2.1
36	DB	55	U	2.1
52	DV	26	ASP	2.1
35	BA	2370	G	2.1
36	BB	21	G	2.1
25	CY	90	LEU	2.1
35	BA	923	C	2.1
35	BA	1053	C	2.1
15	CO	21	ASP	2.1
1	CA	427	U	2.1
49	DS	109	GLY	2.1
1	AA	189	G	2.0
1	CA	1087	G	2.0
35	BA	932	G	2.0
35	BA	1485	G	2.0
9	AI	105	ASP	2.0
11	AK	20	TYR	2.0
16	CP	39	TYR	2.0
49	DS	82	ILE	2.0
1	AA	1165	C	2.0
1	AA	1303	C	2.0
1	CA	1200	C	2.0
1	CA	1321	C	2.0
45	BO	58	VAL	2.0
9	CI	30	GLY	2.0
25	AY	95	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
35	DA	2473	U	2.0
2	AB	190	THR	2.0
12	AL	61	THR	2.0
1	AA	1286	A	2.0
1	CA	451	A	2.0
7	CG	61	VAL	2.0
26	B0	15	ASP	2.0
9	CI	124	GLN	2.0
35	BA	374	A	2.0
35	DA	1545	A	2.0
56	BZ	114	GLY	2.0
35	DA	1485	G	2.0
36	DB	54	G	2.0
16	CP	42	ARG	2.0
35	DA	1043	C	2.0
35	DA	1504	C	2.0
2	CB	121	LEU	2.0
3	AC	67	THR	2.0
7	CG	118	VAL	2.0
1	CA	1090	U	2.0
35	DA	403	U	2.0
35	DA	1175	U	2.0
41	DG	75	LYS	2.0
53	BW	52	GLU	2.0
4	CD	181	MET	2.0
37	BC	190	ARG	2.0
35	BA	300	A	2.0
49	DS	53	SER	2.0
3	CC	188	LEU	2.0
1	CA	812	C	2.0
1	CA	1310	G	2.0
2	CB	43	ASP	2.0
35	DA	219	G	2.0
35	DA	2876	G	2.0
53	BW	112	GLY	2.0
35	BA	218	A	2.0
35	BA	332	A	2.0
2	CB	31	TYR	2.0
7	AG	87	VAL	2.0
44	BN	15	LEU	2.0
1	AA	19	C	2.0
1	AA	255	G	2.0

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Mol	Chain	Res	Type	RSRZ
1	AA	1228	C	2.0
22	CV	26	G	2.0
35	DA	2276	G	2.0
36	BB	3	C	2.0
16	CP	44	THR	2.0
31	B5	51	TYR	2.0
41	DG	82	LEU	2.0
26	D0	25	ARG	2.0
37	DC	159	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
23	5MU	CW	55	21/22	0.21	0.71	177,200,200,200	0
23	5MU	AW	55	21/22	0.84	0.25	130,173,200,200	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
57	MG	BB	220	1/1	-0.38	1.19	71,71,71,71	1
57	MG	CA	1796	1/1	-0.36	0.27	95,95,95,95	1
57	MG	CW	111	1/1	-0.29	0.56	136,136,136,136	1
57	MG	BA	3423	1/1	-0.04	0.63	54,54,54,54	1
57	MG	BA	3231	1/1	0.01	0.83	102,102,102,102	0
57	MG	CW	110	1/1	0.05	2.15	66,66,66,66	1
57	MG	AW	118	1/1	0.12	0.32	69,69,69,69	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3212	1/1	0.12	0.71	86,86,86,86	1
57	MG	CA	1746	1/1	0.13	0.92	74,74,74,74	0
57	MG	DA	3090	1/1	0.13	0.55	97,97,97,97	0
57	MG	DA	3301	1/1	0.18	1.59	55,55,55,55	1
57	MG	CW	116	1/1	0.24	0.73	107,107,107,107	0
57	MG	DA	3439	1/1	0.28	1.57	85,85,85,85	1
57	MG	AA	1789	1/1	0.29	0.45	111,111,111,111	0
57	MG	CA	1759	1/1	0.31	2.40	108,108,108,108	1
57	MG	AW	112	1/1	0.34	0.23	64,64,64,64	1
57	MG	CW	123	1/1	0.42	0.32	1,1,1,1	1
57	MG	DB	202	1/1	0.42	0.67	33,33,33,33	1
57	MG	BA	3281	1/1	0.43	1.96	43,43,43,43	1
57	MG	CA	1785	1/1	0.43	0.43	112,112,112,112	0
57	MG	AW	111	1/1	0.43	0.32	76,76,76,76	1
57	MG	DA	3329	1/1	0.44	1.32	62,62,62,62	1
57	MG	AA	1706	1/1	0.45	0.57	58,58,58,58	1
57	MG	CW	115	1/1	0.46	0.85	93,93,93,93	1
57	MG	AA	1780	1/1	0.46	0.68	95,95,95,95	1
57	MG	CW	106	1/1	0.47	1.15	47,47,47,47	1
57	MG	DA	3361	1/1	0.49	0.51	88,88,88,88	0
57	MG	CA	1681	1/1	0.50	0.48	96,96,96,96	1
57	MG	CW	109	1/1	0.51	0.36	30,30,30,30	1
57	MG	AA	1650	1/1	0.51	1.83	20,20,20,20	1
57	MG	BA	3379	1/1	0.51	1.28	50,50,50,50	1
57	MG	AA	1799	1/1	0.52	0.37	58,58,58,58	1
57	MG	BA	3018	1/1	0.52	0.26	35,35,35,35	0
57	MG	BA	3412	1/1	0.53	0.50	34,34,34,34	1
57	MG	CA	1731	1/1	0.54	0.49	54,54,54,54	0
57	MG	CA	1737	1/1	0.55	0.66	1,1,1,1	1
57	MG	CW	117	1/1	0.55	0.26	57,57,57,57	1
57	MG	CW	113	1/1	0.55	0.68	68,68,68,68	1
57	MG	DA	3295	1/1	0.57	0.23	79,79,79,79	0
57	MG	BA	3251	1/1	0.57	0.52	64,64,64,64	0
57	MG	BA	3392	1/1	0.58	0.19	70,70,70,70	0
57	MG	BA	3035	1/1	0.58	0.40	67,67,67,67	1
57	MG	CA	1715	1/1	0.58	0.20	55,55,55,55	0
57	MG	D1	104	1/1	0.59	0.55	74,74,74,74	1
57	MG	BA	3227	1/1	0.60	1.36	71,71,71,71	1
57	MG	AA	1798	1/1	0.60	0.70	52,52,52,52	1
57	MG	DA	3273	1/1	0.60	0.44	97,97,97,97	1
57	MG	DA	3127	1/1	0.60	0.54	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	CA	1786	1/1	0.60	0.49	39,39,39,39	1
57	MG	DA	3422	1/1	0.62	0.22	62,62,62,62	0
57	MG	DA	3176	1/1	0.62	0.62	77,77,77,77	0
57	MG	DA	3221	1/1	0.62	0.79	79,79,79,79	1
57	MG	DA	3371	1/1	0.62	0.19	35,35,35,35	1
57	MG	CA	1725	1/1	0.62	0.64	43,43,43,43	1
57	MG	AA	1743	1/1	0.62	0.65	74,74,74,74	1
57	MG	DA	3270	1/1	0.63	1.08	47,47,47,47	1
57	MG	DA	3140	1/1	0.63	0.44	25,25,25,25	0
57	MG	BB	213	1/1	0.64	0.78	33,33,33,33	1
57	MG	DA	3278	1/1	0.65	0.33	36,36,36,36	1
57	MG	CA	1686	1/1	0.65	0.46	49,49,49,49	0
57	MG	DB	212	1/1	0.66	0.27	55,55,55,55	0
57	MG	AA	1762	1/1	0.66	0.44	58,58,58,58	0
57	MG	BB	210	1/1	0.66	0.89	68,68,68,68	1
57	MG	BA	3287	1/1	0.67	0.43	45,45,45,45	1
57	MG	AA	1615	1/1	0.67	1.36	47,47,47,47	0
57	MG	DH	201	1/1	0.67	0.53	12,12,12,12	1
57	MG	CA	1770	1/1	0.67	1.18	78,78,78,78	1
57	MG	DA	3281	1/1	0.67	0.48	64,64,64,64	0
57	MG	DB	219	1/1	0.67	0.74	56,56,56,56	1
57	MG	BA	3324	1/1	0.67	0.34	56,56,56,56	0
57	MG	BA	3014	1/1	0.68	0.76	51,51,51,51	0
57	MG	BA	3207	1/1	0.68	0.75	1,1,1,1	1
57	MG	CA	1673	1/1	0.68	0.11	57,57,57,57	0
57	MG	CA	1658	1/1	0.68	0.99	24,24,24,24	1
57	MG	AA	1716	1/1	0.68	0.26	60,60,60,60	0
57	MG	DA	3241	1/1	0.68	0.86	50,50,50,50	0
57	MG	BA	3101	1/1	0.68	0.43	1,1,1,1	1
57	MG	AW	108	1/1	0.69	1.09	24,24,24,24	1
57	MG	DA	3366	1/1	0.69	0.40	46,46,46,46	1
57	MG	AW	123	1/1	0.69	0.47	1,1,1,1	1
57	MG	BB	209	1/1	0.69	0.50	108,108,108,108	0
57	MG	CE	202	1/1	0.69	0.50	85,85,85,85	1
57	MG	CA	1792	1/1	0.70	0.15	16,16,16,16	1
57	MG	AA	1791	1/1	0.70	1.23	35,35,35,35	1
57	MG	BB	217	1/1	0.70	0.23	49,49,49,49	0
57	MG	BA	3334	1/1	0.70	0.44	51,51,51,51	1
57	MG	CW	118	1/1	0.70	0.43	1,1,1,1	1
57	MG	CA	1726	1/1	0.70	0.33	36,36,36,36	0
57	MG	DA	3159	1/1	0.70	0.30	21,21,21,21	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	CA	1623	1/1	0.70	0.34	84,84,84,84	1
57	MG	DA	3101	1/1	0.70	0.43	17,17,17,17	1
57	MG	DA	3434	1/1	0.71	0.38	98,98,98,98	0
57	MG	DA	3215	1/1	0.71	0.58	17,17,17,17	1
57	MG	DA	3299	1/1	0.71	0.51	13,13,13,13	1
57	MG	BA	3153	1/1	0.71	0.35	35,35,35,35	1
57	MG	AV	102	1/1	0.71	0.41	73,73,73,73	0
57	MG	CA	1602	1/1	0.71	0.61	72,72,72,72	0
57	MG	DA	3370	1/1	0.71	0.47	100,100,100,100	1
57	MG	DA	3363	1/1	0.71	0.44	13,13,13,13	1
57	MG	DA	3214	1/1	0.71	0.35	64,64,64,64	1
57	MG	DA	3367	1/1	0.72	0.56	95,95,95,95	1
57	MG	BA	3397	1/1	0.72	0.41	44,44,44,44	0
57	MG	CV	102	1/1	0.72	0.45	57,57,57,57	0
57	MG	BA	3286	1/1	0.73	0.43	32,32,32,32	1
57	MG	DA	3262	1/1	0.73	1.45	79,79,79,79	1
57	MG	CA	1788	1/1	0.73	0.33	1,1,1,1	1
57	MG	CA	1783	1/1	0.73	0.10	33,33,33,33	1
57	MG	CV	101	1/1	0.73	0.15	41,41,41,41	0
57	MG	CA	1765	1/1	0.73	0.64	85,85,85,85	0
57	MG	AA	1749	1/1	0.73	0.45	1,1,1,1	1
57	MG	CA	1738	1/1	0.74	0.93	67,67,67,67	0
57	MG	CN	101	1/1	0.74	0.51	65,65,65,65	0
57	MG	CA	1718	1/1	0.74	0.94	59,59,59,59	0
57	MG	AA	1769	1/1	0.74	1.24	48,48,48,48	0
57	MG	BA	3327	1/1	0.74	0.13	46,46,46,46	1
57	MG	BA	3316	1/1	0.74	0.59	1,1,1,1	1
57	MG	DA	3006	1/1	0.74	0.48	93,93,93,93	0
57	MG	DA	3296	1/1	0.75	0.68	33,33,33,33	1
57	MG	BA	3322	1/1	0.75	0.77	21,21,21,21	1
57	MG	CW	104	1/1	0.75	0.21	35,35,35,35	1
57	MG	DN	202	1/1	0.75	0.29	31,31,31,31	1
57	MG	CA	1707	1/1	0.75	0.34	32,32,32,32	0
57	MG	BA	3002	1/1	0.75	0.39	16,16,16,16	1
57	MG	DA	3326	1/1	0.75	0.85	96,96,96,96	1
57	MG	AA	1653	1/1	0.76	0.79	93,93,93,93	1
57	MG	CA	1709	1/1	0.76	0.19	46,46,46,46	0
57	MG	BB	204	1/1	0.76	0.75	76,76,76,76	1
57	MG	BA	3168	1/1	0.76	0.37	29,29,29,29	0
57	MG	DA	3049	1/1	0.76	0.99	43,43,43,43	0
57	MG	CA	1697	1/1	0.76	0.36	59,59,59,59	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	D2	2601	1/1	0.76	0.47	73,73,73,73	1
57	MG	DA	3319	1/1	0.76	0.25	116,116,116,116	1
57	MG	DA	3358	1/1	0.76	0.71	11,11,11,11	1
57	MG	DA	3161	1/1	0.76	0.31	112,112,112,112	0
57	MG	BA	3309	1/1	0.76	0.66	47,47,47,47	0
57	MG	DA	3174	1/1	0.76	0.56	50,50,50,50	1
57	MG	DA	3346	1/1	0.76	0.43	115,115,115,115	0
57	MG	AW	102	1/1	0.76	0.39	82,82,82,82	1
57	MG	CA	1804	1/1	0.76	0.32	45,45,45,45	0
57	MG	AW	103	1/1	0.76	0.29	1,1,1,1	1
57	MG	CA	1747	1/1	0.76	0.66	19,19,19,19	1
57	MG	BA	3006	1/1	0.77	0.45	77,77,77,77	0
57	MG	BA	3102	1/1	0.77	0.38	21,21,21,21	1
57	MG	DA	3409	1/1	0.77	1.57	83,83,83,83	1
57	MG	AW	113	1/1	0.77	0.27	60,60,60,60	0
57	MG	AV	103	1/1	0.77	0.14	70,70,70,70	0
57	MG	CA	1646	1/1	0.77	0.44	56,56,56,56	0
57	MG	CA	1736	1/1	0.77	0.45	71,71,71,71	0
57	MG	BN	204	1/1	0.77	0.36	20,20,20,20	1
57	MG	CA	1767	1/1	0.77	1.32	94,94,94,94	0
57	MG	DA	3307	1/1	0.77	0.48	43,43,43,43	0
57	MG	CA	1643	1/1	0.77	1.02	73,73,73,73	1
57	MG	CA	1754	1/1	0.77	0.33	42,42,42,42	0
57	MG	BA	3270	1/1	0.77	0.23	6,6,6,6	0
57	MG	CA	1649	1/1	0.78	0.23	14,14,14,14	1
57	MG	AK	201	1/1	0.78	0.44	33,33,33,33	1
57	MG	AA	1734	1/1	0.78	0.79	36,36,36,36	0
57	MG	CA	1760	1/1	0.78	0.44	40,40,40,40	0
57	MG	CA	1758	1/1	0.78	0.39	21,21,21,21	1
57	MG	DA	3242	1/1	0.78	0.18	100,100,100,100	1
57	MG	DA	3004	1/1	0.78	0.83	64,64,64,64	0
57	MG	DA	3152	1/1	0.78	0.66	62,62,62,62	1
57	MG	BA	3229	1/1	0.78	0.21	39,39,39,39	0
57	MG	AA	1788	1/1	0.78	0.30	40,40,40,40	0
57	MG	DA	3321	1/1	0.78	0.36	85,85,85,85	0
57	MG	CW	102	1/1	0.79	0.27	38,38,38,38	1
57	MG	DX	103	1/1	0.79	0.35	48,48,48,48	1
57	MG	CA	1668	1/1	0.79	0.28	43,43,43,43	0
57	MG	BA	3299	1/1	0.79	1.03	26,26,26,26	1
57	MG	BA	3292	1/1	0.79	0.81	27,27,27,27	1
57	MG	DA	3165	1/1	0.79	0.17	47,47,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3135	1/1	0.79	0.86	12,12,12,12	0
57	MG	DA	3139	1/1	0.79	0.59	53,53,53,53	0
57	MG	DA	3271	1/1	0.79	0.51	113,113,113,113	1
57	MG	AA	1733	1/1	0.79	0.43	5,5,5,5	1
57	MG	BA	3289	1/1	0.79	0.37	60,60,60,60	1
57	MG	DA	3220	1/1	0.79	0.77	86,86,86,86	0
57	MG	CW	112	1/1	0.79	0.47	24,24,24,24	1
57	MG	CA	1797	1/1	0.79	1.54	82,82,82,82	1
57	MG	BA	3005	1/1	0.79	0.37	25,25,25,25	0
57	MG	BA	3206	1/1	0.79	0.35	1,1,1,1	1
57	MG	DA	3303	1/1	0.79	0.54	4,4,4,4	0
57	MG	AA	1785	1/1	0.79	0.15	19,19,19,19	0
57	MG	DA	3400	1/1	0.79	0.80	61,61,61,61	1
57	MG	DA	3355	1/1	0.79	0.19	42,42,42,42	1
57	MG	AW	110	1/1	0.80	0.32	30,30,30,30	1
57	MG	AA	1748	1/1	0.80	0.18	73,73,73,73	0
57	MG	DA	3258	1/1	0.80	0.67	23,23,23,23	1
57	MG	BA	3189	1/1	0.80	0.92	87,87,87,87	0
57	MG	DA	3436	1/1	0.80	0.70	41,41,41,41	0
57	MG	DA	3282	1/1	0.80	0.41	1,1,1,1	1
57	MG	DA	3441	1/1	0.80	1.10	44,44,44,44	1
57	MG	AA	1680	1/1	0.80	0.33	53,53,53,53	0
57	MG	AA	1737	1/1	0.80	0.35	30,30,30,30	1
57	MG	AA	1718	1/1	0.80	0.40	1,1,1,1	1
57	MG	AW	114	1/1	0.80	0.16	78,78,78,78	0
57	MG	AW	121	1/1	0.80	0.49	63,63,63,63	0
57	MG	DA	3385	1/1	0.80	0.38	44,44,44,44	0
57	MG	DB	209	1/1	0.80	0.40	96,96,96,96	1
57	MG	AA	1720	1/1	0.80	0.48	91,91,91,91	0
57	MG	BA	3167	1/1	0.80	0.29	50,50,50,50	0
57	MG	AA	1705	1/1	0.80	0.29	38,38,38,38	0
57	MG	CA	1735	1/1	0.80	0.30	1,1,1,1	1
57	MG	DA	3154	1/1	0.81	0.41	83,83,83,83	1
57	MG	BA	3325	1/1	0.81	0.26	30,30,30,30	0
57	MG	AW	106	1/1	0.81	0.54	26,26,26,26	1
57	MG	CG	201	1/1	0.81	0.35	10,10,10,10	1
57	MG	BA	3337	1/1	0.81	0.69	1,1,1,1	1
57	MG	BA	3402	1/1	0.81	0.22	13,13,13,13	1
57	MG	DA	3016	1/1	0.81	0.18	57,57,57,57	0
57	MG	DA	3392	1/1	0.81	0.50	46,46,46,46	0
57	MG	CA	1674	1/1	0.81	0.24	44,44,44,44	1
57	MG	DA	3163	1/1	0.81	0.47	17,17,17,17	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3332	1/1	0.81	0.34	41,41,41,41	1
57	MG	BA	3343	1/1	0.81	0.28	33,33,33,33	0
57	MG	CA	1803	1/1	0.81	0.30	83,83,83,83	0
57	MG	DA	3424	1/1	0.81	0.22	31,31,31,31	1
57	MG	BA	3390	1/1	0.81	0.56	24,24,24,24	0
57	MG	DA	3087	1/1	0.81	0.46	17,17,17,17	1
57	MG	CW	119	1/1	0.82	0.70	57,57,57,57	1
57	MG	DA	3336	1/1	0.82	0.31	60,60,60,60	0
57	MG	AA	1797	1/1	0.82	0.69	9,9,9,9	1
57	MG	DA	3320	1/1	0.82	0.55	75,75,75,75	0
57	MG	BA	3225	1/1	0.82	0.90	67,67,67,67	0
57	MG	DA	3343	1/1	0.82	0.39	1,1,1,1	1
57	MG	BA	3010	1/1	0.82	0.92	17,17,17,17	1
57	MG	AA	1693	1/1	0.82	0.10	80,80,80,80	0
57	MG	CA	1748	1/1	0.82	0.61	97,97,97,97	1
57	MG	DA	3438	1/1	0.82	1.20	48,48,48,48	1
57	MG	CA	1802	1/1	0.82	1.35	39,39,39,39	0
57	MG	DB	208	1/1	0.82	0.17	53,53,53,53	1
57	MG	CA	1801	1/1	0.82	1.27	2,2,2,2	1
57	MG	DA	3279	1/1	0.82	0.46	75,75,75,75	0
57	MG	DA	3360	1/1	0.82	0.09	5,5,5,5	0
57	MG	CA	1793	1/1	0.82	0.64	74,74,74,74	0
57	MG	CW	107	1/1	0.82	0.99	84,84,84,84	0
57	MG	DB	216	1/1	0.82	0.28	54,54,54,54	0
57	MG	BF	304	1/1	0.82	0.36	9,9,9,9	1
57	MG	CA	1655	1/1	0.82	0.55	69,69,69,69	1
57	MG	DB	218	1/1	0.82	0.25	30,30,30,30	1
57	MG	DA	3100	1/1	0.82	0.36	63,63,63,63	1
57	MG	DA	3382	1/1	0.82	0.37	71,71,71,71	0
57	MG	CA	1692	1/1	0.83	0.32	74,74,74,74	1
57	MG	AA	1764	1/1	0.83	0.26	35,35,35,35	0
57	MG	BA	3315	1/1	0.83	0.56	87,87,87,87	0
57	MG	DA	3316	1/1	0.83	0.41	14,14,14,14	1
57	MG	CA	1794	1/1	0.83	0.70	68,68,68,68	0
57	MG	AA	1792	1/1	0.83	0.52	56,56,56,56	0
57	MG	CA	1756	1/1	0.83	0.54	56,56,56,56	0
57	MG	AA	1640	1/1	0.83	0.33	18,18,18,18	0
57	MG	CA	1743	1/1	0.83	0.33	64,64,64,64	0
57	MG	BA	3443	1/1	0.83	0.14	49,49,49,49	0
57	MG	AA	1715	1/1	0.83	0.98	29,29,29,29	1
57	MG	CA	1684	1/1	0.83	0.35	59,59,59,59	0
57	MG	BA	3285	1/1	0.83	0.54	40,40,40,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3141	1/1	0.83	0.25	6,6,6,6	1
57	MG	BA	3365	1/1	0.83	0.35	36,36,36,36	0
57	MG	AA	1683	1/1	0.83	0.43	1,1,1,1	1
57	MG	AA	1679	1/1	0.83	0.46	12,12,12,12	1
57	MG	CW	101	1/1	0.83	1.79	83,83,83,83	1
57	MG	DA	3181	1/1	0.83	0.36	47,47,47,47	1
57	MG	BA	3264	1/1	0.83	1.20	66,66,66,66	0
57	MG	CA	1704	1/1	0.83	0.82	62,62,62,62	0
57	MG	BA	3103	1/1	0.83	0.95	33,33,33,33	1
57	MG	DA	3252	1/1	0.83	0.34	89,89,89,89	0
57	MG	CA	1744	1/1	0.83	0.18	33,33,33,33	0
57	MG	BA	3188	1/1	0.83	0.16	12,12,12,12	0
57	MG	BA	3369	1/1	0.83	0.15	50,50,50,50	0
57	MG	BA	3418	1/1	0.83	0.39	41,41,41,41	1
57	MG	BA	3158	1/1	0.83	0.35	32,32,32,32	0
57	MG	CA	1640	1/1	0.84	0.27	65,65,65,65	1
57	MG	BA	3030	1/1	0.84	0.34	56,56,56,56	1
57	MG	DA	3141	1/1	0.84	0.68	1,1,1,1	1
57	MG	AA	1781	1/1	0.84	0.33	7,7,7,7	1
57	MG	BA	3373	1/1	0.84	0.29	1,1,1,1	1
57	MG	DA	3372	1/1	0.84	0.71	35,35,35,35	1
57	MG	BA	3417	1/1	0.84	0.54	17,17,17,17	1
57	MG	CA	1645	1/1	0.84	0.30	46,46,46,46	0
57	MG	AA	1668	1/1	0.84	0.31	18,18,18,18	0
57	MG	BA	3445	1/1	0.84	0.51	25,25,25,25	1
57	MG	BA	3411	1/1	0.84	0.22	99,99,99,99	1
57	MG	CA	1728	1/1	0.84	0.37	41,41,41,41	1
57	MG	AA	1702	1/1	0.84	0.48	57,57,57,57	0
57	MG	DA	3206	1/1	0.84	0.70	32,32,32,32	1
57	MG	AW	109	1/1	0.84	0.28	37,37,37,37	1
57	MG	BA	3166	1/1	0.84	0.27	66,66,66,66	0
57	MG	AA	1726	1/1	0.84	0.31	10,10,10,10	0
57	MG	BA	3340	1/1	0.84	0.63	22,22,22,22	1
57	MG	DA	3150	1/1	0.84	0.45	39,39,39,39	1
57	MG	BA	3053	1/1	0.84	0.14	1,1,1,1	0
57	MG	BA	3441	1/1	0.84	0.20	6,6,6,6	1
57	MG	DA	3207	1/1	0.85	0.54	36,36,36,36	1
57	MG	DA	3292	1/1	0.85	0.37	45,45,45,45	0
57	MG	AA	1696	1/1	0.85	0.41	45,45,45,45	1
57	MG	AW	116	1/1	0.85	0.30	61,61,61,61	0
57	MG	CA	1651	1/1	0.85	0.36	50,50,50,50	1
57	MG	DA	3136	1/1	0.85	0.27	9,9,9,9	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1666	1/1	0.85	0.41	47,47,47,47	0
57	MG	DA	3415	1/1	0.85	0.40	3,3,3,3	1
57	MG	BA	3330	1/1	0.85	0.14	35,35,35,35	0
57	MG	CA	1657	1/1	0.85	0.44	67,67,67,67	0
57	MG	DA	3325	1/1	0.85	0.26	72,72,72,72	0
57	MG	CA	1764	1/1	0.85	0.08	49,49,49,49	0
57	MG	BB	208	1/1	0.85	0.16	38,38,38,38	1
57	MG	BA	3238	1/1	0.85	0.69	38,38,38,38	1
57	MG	CA	1641	1/1	0.85	0.14	17,17,17,17	0
57	MG	DA	3225	1/1	0.85	0.16	35,35,35,35	0
57	MG	BB	203	1/1	0.85	0.26	60,60,60,60	0
57	MG	AA	1671	1/1	0.85	0.33	47,47,47,47	0
57	MG	DF	305	1/1	0.85	0.30	1,1,1,1	1
57	MG	DA	3275	1/1	0.85	0.89	2,2,2,2	1
57	MG	BA	3374	1/1	0.85	0.95	28,28,28,28	1
57	MG	CA	1782	1/1	0.85	0.12	28,28,28,28	0
57	MG	BA	3362	1/1	0.85	0.28	38,38,38,38	0
57	MG	DA	3390	1/1	0.85	0.19	34,34,34,34	0
57	MG	DA	3431	1/1	0.85	0.12	16,16,16,16	0
57	MG	CA	1778	1/1	0.85	0.46	1,1,1,1	1
57	MG	CV	104	1/1	0.85	0.35	45,45,45,45	1
57	MG	AA	1603	1/1	0.85	0.23	20,20,20,20	0
57	MG	BA	3434	1/1	0.86	0.20	60,60,60,60	0
57	MG	BA	3321	1/1	0.86	0.12	51,51,51,51	0
57	MG	AA	1676	1/1	0.86	0.24	32,32,32,32	0
57	MG	BA	3348	1/1	0.86	1.26	84,84,84,84	0
57	MG	CA	1702	1/1	0.86	0.64	47,47,47,47	0
57	MG	BA	3405	1/1	0.86	0.36	22,22,22,22	0
57	MG	DA	3426	1/1	0.86	0.24	44,44,44,44	0
57	MG	AA	1800	1/1	0.86	0.29	38,38,38,38	0
57	MG	DA	3245	1/1	0.86	0.60	67,67,67,67	0
57	MG	DA	3302	1/1	0.86	0.28	80,80,80,80	1
57	MG	DA	3323	1/1	0.86	0.14	63,63,63,63	1
57	MG	BA	3178	1/1	0.86	0.38	38,38,38,38	0
57	MG	BA	3294	1/1	0.86	0.54	55,55,55,55	0
57	MG	CA	1733	1/1	0.86	0.39	59,59,59,59	0
57	MG	CA	1774	1/1	0.86	0.32	74,74,74,74	0
57	MG	BA	3157	1/1	0.86	0.20	28,28,28,28	0
57	MG	CA	1771	1/1	0.86	0.08	1,1,1,1	1
57	MG	DA	3138	1/1	0.86	0.26	17,17,17,17	0
57	MG	BA	3232	1/1	0.86	0.34	30,30,30,30	0
57	MG	BA	3437	1/1	0.86	0.32	64,64,64,64	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1760	1/1	0.86	0.24	37,37,37,37	0
57	MG	DA	3396	1/1	0.86	0.60	14,14,14,14	0
57	MG	BA	3088	1/1	0.86	0.31	1,1,1,1	1
57	MG	CW	108	1/1	0.86	0.43	28,28,28,28	1
57	MG	CA	1616	1/1	0.86	0.79	64,64,64,64	0
57	MG	DA	3285	1/1	0.86	0.52	13,13,13,13	1
57	MG	DA	3202	1/1	0.86	0.88	7,7,7,7	1
57	MG	DB	201	1/1	0.87	0.23	22,22,22,22	1
57	MG	BA	3274	1/1	0.87	0.25	55,55,55,55	0
57	MG	AA	1723	1/1	0.87	0.22	55,55,55,55	0
57	MG	AA	1712	1/1	0.87	0.31	50,50,50,50	0
57	MG	BA	3250	1/1	0.87	0.16	24,24,24,24	0
57	MG	CA	1761	1/1	0.87	0.22	1,1,1,1	1
57	MG	DA	3322	1/1	0.87	0.19	34,34,34,34	1
57	MG	CA	1690	1/1	0.87	0.60	20,20,20,20	1
57	MG	DB	204	1/1	0.87	0.43	89,89,89,89	1
57	MG	BA	3282	1/1	0.87	0.31	72,72,72,72	0
57	MG	AA	1701	1/1	0.87	0.58	8,8,8,8	1
57	MG	CA	1617	1/1	0.87	0.72	29,29,29,29	0
57	MG	CA	1724	1/1	0.87	0.36	51,51,51,51	0
57	MG	BA	3408	1/1	0.87	0.24	21,21,21,21	0
57	MG	AA	1725	1/1	0.87	0.49	48,48,48,48	0
57	MG	AA	1646	1/1	0.87	0.42	29,29,29,29	0
57	MG	BA	3029	1/1	0.87	0.64	1,1,1,1	1
57	MG	BA	3269	1/1	0.87	0.21	22,22,22,22	0
57	MG	AW	115	1/1	0.87	0.39	19,19,19,19	1
57	MG	DA	3407	1/1	0.87	0.28	9,9,9,9	0
57	MG	AA	1644	1/1	0.87	0.29	1,1,1,1	0
57	MG	B1	102	1/1	0.87	0.22	7,7,7,7	1
57	MG	BA	3306	1/1	0.87	0.86	50,50,50,50	0
57	MG	CA	1742	1/1	0.87	0.70	1,1,1,1	1
57	MG	AA	1742	1/1	0.87	0.14	36,36,36,36	0
57	MG	AA	1758	1/1	0.87	0.20	11,11,11,11	0
57	MG	DA	3317	1/1	0.87	0.40	78,78,78,78	1
57	MG	AA	1741	1/1	0.87	0.30	48,48,48,48	0
57	MG	AA	1682	1/1	0.87	0.23	46,46,46,46	0
57	MG	BA	3129	1/1	0.87	0.31	8,8,8,8	1
57	MG	DA	3192	1/1	0.87	0.61	27,27,27,27	0
57	MG	DA	3289	1/1	0.87	0.53	36,36,36,36	0
57	MG	BA	3422	1/1	0.87	0.20	30,30,30,30	0
57	MG	BA	3258	1/1	0.87	0.17	32,32,32,32	0
57	MG	BA	3279	1/1	0.87	0.25	35,35,35,35	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AV	101	1/1	0.87	0.37	46,46,46,46	0
57	MG	BA	3323	1/1	0.88	0.35	74,74,74,74	0
57	MG	DA	3102	1/1	0.88	0.17	3,3,3,3	0
57	MG	CA	1687	1/1	0.88	0.44	42,42,42,42	0
57	MG	BA	3130	1/1	0.88	0.28	1,1,1,1	1
57	MG	CA	1610	1/1	0.88	0.22	15,15,15,15	1
57	MG	DA	3069	1/1	0.88	0.35	21,21,21,21	0
57	MG	DA	3387	1/1	0.88	0.26	33,33,33,33	1
57	MG	BA	3266	1/1	0.88	0.82	65,65,65,65	0
57	MG	BA	3440	1/1	0.88	0.13	33,33,33,33	0
57	MG	DA	3312	1/1	0.88	0.31	1,1,1,1	0
57	MG	BA	3356	1/1	0.88	0.29	3,3,3,3	1
57	MG	CA	1763	1/1	0.88	0.48	15,15,15,15	0
57	MG	D1	102	1/1	0.88	0.32	51,51,51,51	0
57	MG	AA	1736	1/1	0.88	0.44	20,20,20,20	0
57	MG	AA	1686	1/1	0.88	0.24	61,61,61,61	0
57	MG	CA	1650	1/1	0.88	0.25	98,98,98,98	1
57	MG	BA	3137	1/1	0.88	0.72	7,7,7,7	0
57	MG	DA	3440	1/1	0.88	0.41	80,80,80,80	0
57	MG	DG	201	1/1	0.88	0.21	38,38,38,38	0
57	MG	CX	103	1/1	0.88	0.28	45,45,45,45	1
57	MG	DA	3318	1/1	0.88	0.63	7,7,7,7	1
57	MG	D3	101	1/1	0.88	0.27	42,42,42,42	0
57	MG	BA	3120	1/1	0.88	0.43	38,38,38,38	0
57	MG	BA	3230	1/1	0.88	0.22	22,22,22,22	0
57	MG	CW	121	1/1	0.88	0.37	28,28,28,28	1
57	MG	DA	3166	1/1	0.88	0.35	43,43,43,43	0
57	MG	CA	1745	1/1	0.88	0.21	112,112,112,112	1
57	MG	BA	3058	1/1	0.88	0.45	3,3,3,3	1
57	MG	BA	3439	1/1	0.88	0.37	25,25,25,25	0
57	MG	DA	3048	1/1	0.88	0.84	56,56,56,56	1
57	MG	DS	201	1/1	0.88	0.21	1,1,1,1	1
57	MG	BA	3204	1/1	0.88	0.26	38,38,38,38	0
57	MG	DA	3236	1/1	0.88	0.40	14,14,14,14	0
57	MG	DA	3042	1/1	0.88	0.22	17,17,17,17	1
57	MG	DA	3328	1/1	0.88	0.63	63,63,63,63	0
57	MG	BA	3128	1/1	0.88	0.28	1,1,1,1	0
57	MG	BA	3276	1/1	0.88	0.82	34,34,34,34	1
57	MG	AW	119	1/1	0.88	1.01	71,71,71,71	1
57	MG	BA	3241	1/1	0.88	0.51	27,27,27,27	0
57	MG	BA	3351	1/1	0.88	0.36	68,68,68,68	1
57	MG	BB	206	1/1	0.88	0.69	13,13,13,13	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3203	1/1	0.88	0.27	10,10,10,10	0
57	MG	DA	3406	1/1	0.88	0.82	57,57,57,57	0
57	MG	BB	215	1/1	0.89	0.25	49,49,49,49	1
57	MG	BA	3284	1/1	0.89	0.29	12,12,12,12	1
57	MG	DA	3022	1/1	0.89	0.26	28,28,28,28	1
57	MG	AX	101	1/1	0.89	0.28	15,15,15,15	1
57	MG	BA	3409	1/1	0.89	0.40	5,5,5,5	0
57	MG	BA	3154	1/1	0.89	0.19	56,56,56,56	0
57	MG	DA	3435	1/1	0.89	0.25	48,48,48,48	0
57	MG	CA	1716	1/1	0.89	0.81	75,75,75,75	0
57	MG	BA	3395	1/1	0.89	0.28	1,1,1,1	1
57	MG	BF	303	1/1	0.89	0.18	36,36,36,36	0
57	MG	BA	3176	1/1	0.89	0.57	21,21,21,21	1
57	MG	CA	1615	1/1	0.89	0.23	42,42,42,42	0
57	MG	AW	117	1/1	0.89	0.20	19,19,19,19	1
57	MG	CW	103	1/1	0.89	0.76	61,61,61,61	1
57	MG	AA	1694	1/1	0.89	0.24	13,13,13,13	1
57	MG	BA	3263	1/1	0.89	0.28	29,29,29,29	0
57	MG	DA	3194	1/1	0.89	0.40	57,57,57,57	0
57	MG	AA	1754	1/1	0.89	0.29	49,49,49,49	1
57	MG	AA	1669	1/1	0.89	0.42	5,5,5,5	0
57	MG	CA	1711	1/1	0.89	0.44	37,37,37,37	1
57	MG	DA	3268	1/1	0.89	0.22	24,24,24,24	0
57	MG	BA	3156	1/1	0.89	0.63	12,12,12,12	1
57	MG	DA	3290	1/1	0.89	0.49	21,21,21,21	0
57	MG	AA	1625	1/1	0.89	0.43	31,31,31,31	0
57	MG	DA	3018	1/1	0.89	0.39	12,12,12,12	0
57	MG	BA	3372	1/1	0.89	0.34	20,20,20,20	1
57	MG	BA	3320	1/1	0.89	0.23	37,37,37,37	1
57	MG	AA	1738	1/1	0.89	0.42	18,18,18,18	1
57	MG	AA	1665	1/1	0.89	0.44	21,21,21,21	0
57	MG	BA	3163	1/1	0.89	0.21	33,33,33,33	0
57	MG	BA	3060	1/1	0.89	0.23	42,42,42,42	0
57	MG	AE	202	1/1	0.89	0.21	37,37,37,37	0
57	MG	BA	3388	1/1	0.89	0.56	4,4,4,4	1
57	MG	AA	1687	1/1	0.89	0.14	61,61,61,61	0
57	MG	CA	1680	1/1	0.89	0.23	75,75,75,75	1
57	MG	DA	3331	1/1	0.89	0.22	79,79,79,79	1
57	MG	CA	1729	1/1	0.89	0.57	63,63,63,63	0
57	MG	AA	1613	1/1	0.89	0.49	11,11,11,11	1
57	MG	BA	3150	1/1	0.89	0.25	22,22,22,22	0
57	MG	BA	3375	1/1	0.89	0.13	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3216	1/1	0.89	0.09	49,49,49,49	0
57	MG	CA	1808	1/1	0.89	0.40	43,43,43,43	1
57	MG	DP	202	1/1	0.89	0.32	1,1,1,1	1
57	MG	BA	3416	1/1	0.90	0.35	13,13,13,13	0
57	MG	BA	3368	1/1	0.90	0.20	1,1,1,1	1
57	MG	AA	1688	1/1	0.90	0.44	12,12,12,12	1
57	MG	AA	1707	1/1	0.90	0.52	53,53,53,53	0
57	MG	AW	122	1/1	0.90	0.34	47,47,47,47	0
57	MG	DB	213	1/1	0.90	0.16	28,28,28,28	1
57	MG	AA	1735	1/1	0.90	0.24	15,15,15,15	0
57	MG	CA	1622	1/1	0.90	0.23	39,39,39,39	0
57	MG	DA	3224	1/1	0.90	0.24	65,65,65,65	1
57	MG	BA	3367	1/1	0.90	0.44	37,37,37,37	0
57	MG	AA	1773	1/1	0.90	0.81	16,16,16,16	1
57	MG	AA	1663	1/1	0.90	0.58	36,36,36,36	0
57	MG	CA	1659	1/1	0.90	0.15	34,34,34,34	1
57	MG	BA	3191	1/1	0.90	0.19	23,23,23,23	0
57	MG	AA	1703	1/1	0.90	0.12	22,22,22,22	0
57	MG	AA	1782	1/1	0.90	0.29	47,47,47,47	0
57	MG	DB	215	1/1	0.90	0.34	63,63,63,63	1
57	MG	BA	3219	1/1	0.90	0.63	1,1,1,1	1
57	MG	BA	3346	1/1	0.90	0.38	43,43,43,43	0
57	MG	DU	203	1/1	0.90	0.26	49,49,49,49	1
57	MG	AA	1752	1/1	0.90	0.29	42,42,42,42	1
57	MG	DP	201	1/1	0.90	0.68	1,1,1,1	1
57	MG	DA	3339	1/1	0.90	0.50	1,1,1,1	1
57	MG	DA	3177	1/1	0.90	0.49	22,22,22,22	0
57	MG	DA	3423	1/1	0.90	0.23	49,49,49,49	0
57	MG	AA	1659	1/1	0.90	0.53	27,27,27,27	0
57	MG	AA	1777	1/1	0.90	0.13	6,6,6,6	0
57	MG	DA	3362	1/1	0.90	0.27	15,15,15,15	0
57	MG	AA	1609	1/1	0.90	0.33	7,7,7,7	1
57	MG	BA	3280	1/1	0.90	0.22	72,72,72,72	0
57	MG	CA	1706	1/1	0.90	0.42	47,47,47,47	1
57	MG	BA	3152	1/1	0.90	0.23	62,62,62,62	0
57	MG	AA	1698	1/1	0.90	0.15	1,1,1,1	0
57	MG	CA	1694	1/1	0.90	0.13	59,59,59,59	0
57	MG	CA	1806	1/1	0.90	0.29	35,35,35,35	0
57	MG	AA	1631	1/1	0.90	0.22	1,1,1,1	1
57	MG	BA	3435	1/1	0.90	0.39	85,85,85,85	0
57	MG	DA	3402	1/1	0.90	0.24	45,45,45,45	1
57	MG	BB	218	1/1	0.90	0.11	18,18,18,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3407	1/1	0.90	0.40	10,10,10,10	0
57	MG	AA	1767	1/1	0.90	0.27	21,21,21,21	0
57	MG	AA	1645	1/1	0.90	0.22	14,14,14,14	1
57	MG	BA	3297	1/1	0.90	0.26	12,12,12,12	0
57	MG	CA	1631	1/1	0.90	0.23	57,57,57,57	0
57	MG	BA	3328	1/1	0.90	0.13	33,33,33,33	0
57	MG	BA	3213	1/1	0.90	0.42	49,49,49,49	1
57	MG	AD	302	1/1	0.90	0.44	28,28,28,28	0
57	MG	AA	1765	1/1	0.90	0.13	41,41,41,41	0
57	MG	CW	105	1/1	0.90	0.99	79,79,79,79	1
57	MG	DA	3369	1/1	0.90	0.45	142,142,142,142	1
57	MG	CA	1739	1/1	0.90	0.27	32,32,32,32	0
57	MG	AA	1746	1/1	0.91	0.39	29,29,29,29	0
57	MG	BA	3252	1/1	0.91	0.31	23,23,23,23	0
57	MG	DA	3026	1/1	0.91	0.67	1,1,1,1	0
57	MG	DA	3086	1/1	0.91	0.45	1,1,1,1	1
57	MG	BA	3145	1/1	0.91	0.29	4,4,4,4	0
57	MG	BA	3100	1/1	0.91	0.33	1,1,1,1	1
57	MG	DA	3344	1/1	0.91	0.15	20,20,20,20	0
57	MG	DA	3269	1/1	0.91	0.52	57,57,57,57	1
57	MG	BA	3331	1/1	0.91	0.14	11,11,11,11	1
57	MG	BA	3020	1/1	0.91	0.24	21,21,21,21	0
57	MG	CA	1741	1/1	0.91	0.22	1,1,1,1	1
57	MG	DA	3330	1/1	0.91	0.30	73,73,73,73	0
57	MG	BA	3406	1/1	0.91	0.45	19,19,19,19	0
57	MG	BA	3335	1/1	0.91	0.36	1,1,1,1	1
57	MG	BA	3107	1/1	0.91	0.17	23,23,23,23	0
57	MG	BA	3112	1/1	0.91	0.23	34,34,34,34	1
57	MG	DE	302	1/1	0.91	0.37	50,50,50,50	0
57	MG	CL	202	1/1	0.91	0.33	5,5,5,5	1
57	MG	DA	3227	1/1	0.91	0.43	73,73,73,73	0
57	MG	AA	1674	1/1	0.91	0.21	30,30,30,30	1
57	MG	DA	3144	1/1	0.91	0.21	1,1,1,1	0
57	MG	DB	206	1/1	0.91	0.22	72,72,72,72	0
57	MG	BA	3333	1/1	0.91	0.10	41,41,41,41	0
57	MG	BA	3224	1/1	0.91	1.25	80,80,80,80	0
57	MG	DA	3149	1/1	0.91	0.69	27,27,27,27	0
57	MG	BA	3182	1/1	0.91	0.30	13,13,13,13	0
57	MG	CA	1798	1/1	0.91	0.37	43,43,43,43	1
57	MG	CA	1768	1/1	0.91	0.59	5,5,5,5	1
57	MG	CA	1653	1/1	0.91	0.11	41,41,41,41	0
57	MG	DA	3395	1/1	0.91	0.64	1,1,1,1	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3310	1/1	0.91	0.86	7,7,7,7	1
57	MG	AA	1724	1/1	0.91	0.22	34,34,34,34	1
57	MG	CA	1664	1/1	0.91	0.10	54,54,54,54	1
57	MG	DA	3195	1/1	0.91	0.16	15,15,15,15	0
57	MG	CA	1727	1/1	0.91	1.14	92,92,92,92	0
57	MG	CA	1772	1/1	0.91	0.18	60,60,60,60	0
57	MG	BA	3389	1/1	0.91	0.65	31,31,31,31	0
57	MG	BA	3222	1/1	0.91	0.14	1,1,1,1	0
57	MG	DA	3218	1/1	0.91	0.42	12,12,12,12	0
57	MG	BF	302	1/1	0.91	0.76	15,15,15,15	1
57	MG	CA	1672	1/1	0.91	0.27	44,44,44,44	0
57	MG	BA	3431	1/1	0.91	0.58	12,12,12,12	0
57	MG	BA	3391	1/1	0.91	0.33	30,30,30,30	0
57	MG	BA	3370	1/1	0.91	0.31	37,37,37,37	0
57	MG	BA	3332	1/1	0.91	0.37	64,64,64,64	0
57	MG	BA	3415	1/1	0.91	0.17	18,18,18,18	1
57	MG	BA	3257	1/1	0.91	0.23	42,42,42,42	0
57	MG	AA	1755	1/1	0.91	0.25	47,47,47,47	0
57	MG	DA	3324	1/1	0.91	0.15	40,40,40,40	0
57	MG	BA	3338	1/1	0.92	0.17	46,46,46,46	0
57	MG	CA	1675	1/1	0.92	0.84	47,47,47,47	1
57	MG	DA	3037	1/1	0.92	0.17	1,1,1,1	0
57	MG	DA	3354	1/1	0.92	0.24	6,6,6,6	1
57	MG	DA	3338	1/1	0.92	0.29	16,16,16,16	1
57	MG	CA	1683	1/1	0.92	0.45	19,19,19,19	0
57	MG	AA	1794	1/1	0.92	0.29	40,40,40,40	0
57	MG	CA	1612	1/1	0.92	0.28	75,75,75,75	0
57	MG	AE	201	1/1	0.92	0.19	28,28,28,28	0
57	MG	DA	3012	1/1	0.92	0.41	1,1,1,1	0
57	MG	DA	3057	1/1	0.92	0.19	3,3,3,3	0
57	MG	DA	3368	1/1	0.92	0.36	78,78,78,78	0
57	MG	BN	202	1/1	0.92	0.62	3,3,3,3	1
57	MG	BA	3380	1/1	0.92	0.17	63,63,63,63	0
57	MG	AA	1611	1/1	0.92	0.12	24,24,24,24	0
57	MG	DA	3403	1/1	0.92	0.47	47,47,47,47	1
57	MG	DA	3414	1/1	0.92	0.18	16,16,16,16	0
57	MG	BA	3001	1/1	0.92	0.38	8,8,8,8	0
57	MG	AA	1616	1/1	0.92	0.12	24,24,24,24	1
57	MG	BA	3209	1/1	0.92	0.08	18,18,18,18	0
57	MG	BA	3275	1/1	0.92	0.14	37,37,37,37	0
57	MG	BA	3245	1/1	0.92	0.09	52,52,52,52	1
57	MG	AA	1704	1/1	0.92	0.20	11,11,11,11	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1803	1/1	0.92	0.24	9,9,9,9	0
57	MG	AA	1761	1/1	0.92	0.19	13,13,13,13	0
57	MG	CA	1666	1/1	0.92	0.28	1,1,1,1	0
57	MG	BA	3179	1/1	0.92	0.18	30,30,30,30	0
57	MG	CA	1700	1/1	0.92	0.25	10,10,10,10	0
57	MG	DA	3107	1/1	0.92	0.38	34,34,34,34	0
57	MG	DA	3340	1/1	0.92	0.15	59,59,59,59	0
57	MG	AA	1655	1/1	0.92	0.33	31,31,31,31	0
57	MG	AA	1602	1/1	0.92	0.15	64,64,64,64	0
57	MG	AA	1721	1/1	0.92	0.43	44,44,44,44	0
57	MG	DA	3264	1/1	0.92	0.24	25,25,25,25	0
57	MG	BA	3038	1/1	0.92	0.34	22,22,22,22	1
57	MG	DA	3420	1/1	0.92	0.26	14,14,14,14	1
57	MG	BA	3143	1/1	0.92	0.45	1,1,1,1	1
57	MG	BA	3267	1/1	0.92	0.24	34,34,34,34	0
57	MG	BA	3353	1/1	0.92	0.54	41,41,41,41	1
57	MG	AA	1608	1/1	0.92	0.06	3,3,3,3	0
57	MG	AD	301	1/1	0.92	0.21	15,15,15,15	1
57	MG	BA	3132	1/1	0.92	0.12	24,24,24,24	1
57	MG	CA	1712	1/1	0.92	0.50	19,19,19,19	0
57	MG	DA	3005	1/1	0.92	0.11	22,22,22,22	0
57	MG	BA	3117	1/1	0.92	0.18	8,8,8,8	0
57	MG	AA	1790	1/1	0.92	0.28	35,35,35,35	0
57	MG	BN	203	1/1	0.92	0.17	56,56,56,56	1
57	MG	CA	1721	1/1	0.92	0.20	17,17,17,17	1
57	MG	BA	3125	1/1	0.92	0.45	1,1,1,1	0
57	MG	BA	3009	1/1	0.92	0.35	1,1,1,1	0
57	MG	AA	1651	1/1	0.92	0.16	32,32,32,32	0
57	MG	CA	1730	1/1	0.92	0.23	1,1,1,1	0
57	MG	BG	201	1/1	0.92	0.09	17,17,17,17	0
57	MG	AA	1787	1/1	0.92	0.19	53,53,53,53	1
57	MG	BA	3239	1/1	0.92	0.20	14,14,14,14	1
57	MG	AA	1623	1/1	0.92	0.23	16,16,16,16	0
57	MG	DA	3345	1/1	0.92	0.15	2,2,2,2	1
57	MG	AA	1753	1/1	0.92	0.41	48,48,48,48	0
57	MG	AA	1634	1/1	0.92	0.11	28,28,28,28	0
57	MG	BA	3197	1/1	0.92	0.20	25,25,25,25	1
57	MG	BA	3308	1/1	0.92	0.97	30,30,30,30	1
57	MG	AA	1621	1/1	0.92	0.12	3,3,3,3	0
57	MG	BA	3421	1/1	0.92	0.69	39,39,39,39	1
57	MG	CA	1607	1/1	0.92	0.24	54,54,54,54	0
57	MG	BA	3387	1/1	0.92	0.15	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1739	1/1	0.92	0.44	44,44,44,44	0
57	MG	BA	3429	1/1	0.92	0.16	32,32,32,32	0
57	MG	CA	1632	1/1	0.92	0.22	36,36,36,36	0
57	MG	CA	1713	1/1	0.92	0.25	40,40,40,40	1
57	MG	DB	205	1/1	0.92	0.40	5,5,5,5	1
57	MG	DA	3126	1/1	0.92	0.20	1,1,1,1	0
57	MG	CA	1613	1/1	0.92	0.70	10,10,10,10	1
57	MG	BA	3169	1/1	0.92	0.33	33,33,33,33	0
57	MG	DA	3389	1/1	0.92	0.26	33,33,33,33	1
57	MG	DA	3309	1/1	0.92	0.34	44,44,44,44	0
57	MG	CA	1719	1/1	0.93	0.20	56,56,56,56	1
57	MG	DA	3008	1/1	0.93	0.41	1,1,1,1	0
57	MG	BA	3123	1/1	0.93	0.58	35,35,35,35	0
57	MG	BA	3272	1/1	0.93	0.15	1,1,1,1	1
57	MG	DA	3286	1/1	0.93	0.40	12,12,12,12	0
57	MG	BA	3091	1/1	0.93	0.17	1,1,1,1	0
57	MG	BA	3295	1/1	0.93	0.12	1,1,1,1	0
57	MG	DA	3377	1/1	0.93	0.84	62,62,62,62	1
57	MG	AA	1728	1/1	0.93	0.15	34,34,34,34	0
57	MG	DA	3180	1/1	0.93	0.39	47,47,47,47	0
57	MG	AA	1629	1/1	0.93	0.10	22,22,22,22	0
57	MG	AA	1656	1/1	0.93	0.26	31,31,31,31	0
57	MG	AA	1648	1/1	0.93	0.16	3,3,3,3	1
57	MG	CA	1644	1/1	0.93	0.24	31,31,31,31	0
57	MG	BA	3359	1/1	0.93	0.39	21,21,21,21	1
57	MG	DA	3014	1/1	0.93	0.24	14,14,14,14	0
57	MG	DA	3432	1/1	0.93	0.41	50,50,50,50	0
57	MG	CA	1769	1/1	0.93	0.23	28,28,28,28	0
57	MG	CA	1639	1/1	0.93	0.44	34,34,34,34	0
57	MG	BA	3383	1/1	0.93	0.96	39,39,39,39	1
57	MG	DA	3428	1/1	0.93	0.55	17,17,17,17	0
57	MG	BA	3386	1/1	0.93	0.14	13,13,13,13	0
57	MG	CA	1660	1/1	0.93	0.16	8,8,8,8	0
57	MG	BA	3364	1/1	0.93	0.19	12,12,12,12	0
57	MG	BA	3089	1/1	0.93	0.56	6,6,6,6	1
57	MG	CW	120	1/1	0.93	0.27	20,20,20,20	1
57	MG	AA	1717	1/1	0.93	0.19	12,12,12,12	1
57	MG	BA	3339	1/1	0.93	0.39	3,3,3,3	0
57	MG	AA	1719	1/1	0.93	0.26	1,1,1,1	1
57	MG	BA	3290	1/1	0.93	1.13	33,33,33,33	1
57	MG	CA	1757	1/1	0.93	0.65	35,35,35,35	0
57	MG	DA	3223	1/1	0.93	0.61	12,12,12,12	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3345	1/1	0.93	0.84	1,1,1,1	1
57	MG	AA	1695	1/1	0.93	0.45	42,42,42,42	0
57	MG	BA	3401	1/1	0.93	0.68	42,42,42,42	0
57	MG	DA	3219	1/1	0.93	0.70	21,21,21,21	1
57	MG	BA	3086	1/1	0.93	0.52	1,1,1,1	0
57	MG	CA	1775	1/1	0.93	0.22	1,1,1,1	1
57	MG	BA	3404	1/1	0.93	0.19	108,108,108,108	0
57	MG	CA	1629	1/1	0.93	0.11	2,2,2,2	1
57	MG	DA	3028	1/1	0.93	0.11	59,59,59,59	1
57	MG	BA	3385	1/1	0.93	0.16	7,7,7,7	0
57	MG	DA	3373	1/1	0.93	0.26	74,74,74,74	0
57	MG	CW	114	1/1	0.93	0.17	4,4,4,4	1
57	MG	CA	1679	1/1	0.93	0.21	22,22,22,22	0
57	MG	BA	3205	1/1	0.93	0.15	23,23,23,23	0
57	MG	DA	3265	1/1	0.93	0.52	70,70,70,70	0
57	MG	BA	3184	1/1	0.93	0.11	44,44,44,44	0
57	MG	DA	3419	1/1	0.93	0.59	58,58,58,58	0
57	MG	AA	1802	1/1	0.93	0.21	18,18,18,18	0
57	MG	BA	3228	1/1	0.93	0.24	50,50,50,50	0
57	MG	BA	3253	1/1	0.93	0.38	56,56,56,56	1
57	MG	BA	3159	1/1	0.93	0.40	1,1,1,1	0
57	MG	BB	214	1/1	0.93	0.15	34,34,34,34	0
57	MG	DA	3201	1/1	0.93	0.14	15,15,15,15	0
57	MG	BA	3140	1/1	0.93	0.24	24,24,24,24	1
57	MG	AA	1684	1/1	0.93	0.11	25,25,25,25	0
57	MG	BA	3350	1/1	0.93	0.35	70,70,70,70	0
57	MG	BA	3246	1/1	0.93	0.30	23,23,23,23	1
57	MG	DA	3045	1/1	0.93	0.21	10,10,10,10	0
57	MG	AA	1710	1/1	0.93	0.38	113,113,113,113	1
57	MG	DA	3071	1/1	0.93	0.38	1,1,1,1	0
57	MG	BA	3093	1/1	0.93	0.14	1,1,1,1	0
57	MG	BA	3349	1/1	0.93	0.34	86,86,86,86	0
57	MG	DA	3147	1/1	0.93	0.21	22,22,22,22	0
57	MG	CA	1784	1/1	0.93	0.23	1,1,1,1	1
57	MG	DA	3231	1/1	0.93	0.43	14,14,14,14	0
57	MG	DA	3170	1/1	0.93	0.33	26,26,26,26	1
57	MG	CA	1652	1/1	0.93	0.49	1,1,1,1	1
57	MG	DA	3145	1/1	0.93	0.46	56,56,56,56	1
57	MG	BA	3336	1/1	0.93	0.19	8,8,8,8	1
57	MG	BA	3303	1/1	0.93	0.40	8,8,8,8	1
57	MG	CA	1654	1/1	0.93	0.24	1,1,1,1	0
57	MG	DA	3388	1/1	0.93	0.74	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3376	1/1	0.93	0.22	1,1,1,1	0
57	MG	BA	3277	1/1	0.93	0.21	28,28,28,28	1
57	MG	DA	3276	1/1	0.93	0.23	47,47,47,47	0
57	MG	DA	3337	1/1	0.93	0.32	92,92,92,92	0
57	MG	DA	3233	1/1	0.93	0.29	1,1,1,1	1
57	MG	DA	3405	1/1	0.93	0.36	30,30,30,30	1
57	MG	DA	3156	1/1	0.93	0.21	21,21,21,21	0
57	MG	BA	3425	1/1	0.93	0.12	23,23,23,23	0
57	MG	DA	3179	1/1	0.93	0.40	52,52,52,52	1
57	MG	BA	3366	1/1	0.93	0.52	38,38,38,38	1
57	MG	CA	1717	1/1	0.93	0.15	1,1,1,1	0
57	MG	BA	3310	1/1	0.93	0.26	5,5,5,5	1
57	MG	CA	1780	1/1	0.93	0.72	31,31,31,31	1
57	MG	BA	3393	1/1	0.93	0.22	14,14,14,14	0
57	MG	CA	1693	1/1	0.93	0.08	14,14,14,14	1
57	MG	BU	201	1/1	0.93	0.24	12,12,12,12	1
57	MG	AW	120	1/1	0.93	0.32	58,58,58,58	1
57	MG	DA	3084	1/1	0.93	0.22	13,13,13,13	0
57	MG	BX	102	1/1	0.93	0.18	1,1,1,1	1
57	MG	CW	122	1/1	0.93	0.22	70,70,70,70	0
57	MG	BA	3139	1/1	0.93	0.38	8,8,8,8	1
57	MG	BA	3198	1/1	0.94	0.26	4,4,4,4	0
57	MG	DA	3274	1/1	0.94	0.05	65,65,65,65	0
57	MG	BA	3214	1/1	0.94	0.16	106,106,106,106	1
57	MG	DA	3046	1/1	0.94	0.41	1,1,1,1	0
57	MG	CA	1698	1/1	0.94	0.11	22,22,22,22	0
57	MG	BA	3341	1/1	0.94	0.16	4,4,4,4	1
57	MG	BA	3248	1/1	0.94	0.18	10,10,10,10	0
57	MG	DA	3253	1/1	0.94	0.52	25,25,25,25	0
57	MG	BA	3399	1/1	0.94	0.22	56,56,56,56	1
57	MG	DA	3010	1/1	0.94	0.17	1,1,1,1	0
57	MG	DA	3228	1/1	0.94	0.56	41,41,41,41	0
57	MG	DA	3248	1/1	0.94	0.33	5,5,5,5	1
57	MG	BA	3318	1/1	0.94	0.32	1,1,1,1	0
57	MG	BA	3073	1/1	0.94	0.47	1,1,1,1	0
57	MG	BA	3043	1/1	0.94	0.19	26,26,26,26	0
57	MG	BA	3442	1/1	0.94	0.16	33,33,33,33	0
57	MG	AA	1757	1/1	0.94	0.33	51,51,51,51	1
57	MG	BA	3135	1/1	0.94	0.17	38,38,38,38	0
57	MG	AA	1766	1/1	0.94	0.19	1,1,1,1	1
57	MG	CA	1620	1/1	0.94	0.13	34,34,34,34	0
57	MG	DA	3247	1/1	0.94	0.30	13,13,13,13	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3026	1/1	0.94	0.16	8,8,8,8	0
57	MG	DA	3430	1/1	0.94	0.21	13,13,13,13	0
57	MG	BA	3396	1/1	0.94	0.56	15,15,15,15	0
57	MG	DD	303	1/1	0.94	0.20	1,1,1,1	0
57	MG	DA	3115	1/1	0.94	0.41	6,6,6,6	0
57	MG	AA	1747	1/1	0.94	0.27	5,5,5,5	0
57	MG	AA	1610	1/1	0.94	0.19	29,29,29,29	0
57	MG	BA	3144	1/1	0.94	0.18	40,40,40,40	0
57	MG	AA	1639	1/1	0.94	0.29	8,8,8,8	1
57	MG	BA	3118	1/1	0.94	0.24	33,33,33,33	1
57	MG	DA	3164	1/1	0.94	0.09	72,72,72,72	0
57	MG	CA	1805	1/1	0.94	0.38	1,1,1,1	0
57	MG	DA	3226	1/1	0.94	0.34	48,48,48,48	0
57	MG	AA	1744	1/1	0.94	0.42	22,22,22,22	0
57	MG	CA	1776	1/1	0.94	0.21	32,32,32,32	0
57	MG	BA	3394	1/1	0.94	0.36	7,7,7,7	1
57	MG	DA	3210	1/1	0.94	0.12	93,93,93,93	1
57	MG	DA	3187	1/1	0.94	0.72	83,83,83,83	1
57	MG	BB	202	1/1	0.94	0.18	9,9,9,9	1
57	MG	CA	1720	1/1	0.94	0.14	4,4,4,4	1
57	MG	DA	3359	1/1	0.94	0.17	40,40,40,40	0
57	MG	AA	1667	1/1	0.94	0.38	17,17,17,17	0
57	MG	DB	207	1/1	0.94	0.26	16,16,16,16	1
57	MG	BA	3109	1/1	0.94	0.45	3,3,3,3	0
57	MG	AA	1763	1/1	0.94	0.64	40,40,40,40	1
57	MG	DA	3349	1/1	0.94	0.21	10,10,10,10	1
57	MG	AA	1779	1/1	0.94	0.26	1,1,1,1	1
57	MG	AA	1713	1/1	0.94	0.14	25,25,25,25	0
57	MG	BA	3307	1/1	0.94	0.36	35,35,35,35	0
57	MG	CE	201	1/1	0.94	0.66	78,78,78,78	0
57	MG	AA	1652	1/1	0.94	0.44	11,11,11,11	0
57	MG	BA	3215	1/1	0.94	0.72	38,38,38,38	1
57	MG	BA	3180	1/1	0.94	0.45	1,1,1,1	1
57	MG	AA	1643	1/1	0.94	0.24	43,43,43,43	0
57	MG	AA	1691	1/1	0.94	0.28	69,69,69,69	0
57	MG	CA	1703	1/1	0.94	0.27	12,12,12,12	0
57	MG	BA	3164	1/1	0.94	0.45	56,56,56,56	0
57	MG	AA	1740	1/1	0.94	0.13	35,35,35,35	0
57	MG	B5	102	1/1	0.94	0.36	48,48,48,48	0
57	MG	DA	3341	1/1	0.94	0.55	20,20,20,20	0
57	MG	DA	3365	1/1	0.94	0.21	24,24,24,24	1
57	MG	DA	3357	1/1	0.94	0.07	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1732	1/1	0.94	0.28	31,31,31,31	0
57	MG	CA	1656	1/1	0.94	0.16	19,19,19,19	1
57	MG	BA	3271	1/1	0.94	0.21	9,9,9,9	0
57	MG	BA	3347	1/1	0.94	0.50	1,1,1,1	0
57	MG	DA	3217	1/1	0.94	0.25	31,31,31,31	0
57	MG	BA	3175	1/1	0.94	0.12	21,21,21,21	0
57	MG	BA	3360	1/1	0.94	0.27	10,10,10,10	0
57	MG	BA	3291	1/1	0.94	0.15	7,7,7,7	0
57	MG	DA	3033	1/1	0.94	0.20	13,13,13,13	0
57	MG	CA	1677	1/1	0.94	0.20	36,36,36,36	0
57	MG	DA	3103	1/1	0.94	0.30	1,1,1,1	0
57	MG	AA	1793	1/1	0.94	0.19	6,6,6,6	1
57	MG	BA	3413	1/1	0.94	0.37	21,21,21,21	1
57	MG	DA	3277	1/1	0.94	0.17	22,22,22,22	0
57	MG	DA	3238	1/1	0.94	0.40	1,1,1,1	0
57	MG	DB	210	1/1	0.94	0.22	49,49,49,49	1
57	MG	AA	1759	1/1	0.94	0.14	20,20,20,20	0
57	MG	CA	1762	1/1	0.94	0.29	7,7,7,7	1
57	MG	BB	219	1/1	0.94	0.45	73,73,73,73	1
57	MG	BA	3201	1/1	0.94	0.20	31,31,31,31	0
57	MG	BA	3011	1/1	0.94	0.43	15,15,15,15	0
57	MG	BO	201	1/1	0.94	0.42	1,1,1,1	1
57	MG	DA	3189	1/1	0.94	0.17	59,59,59,59	0
57	MG	BA	3040	1/1	0.94	0.09	16,16,16,16	0
57	MG	BA	3099	1/1	0.94	0.28	45,45,45,45	1
57	MG	BA	3094	1/1	0.94	0.15	25,25,25,25	1
57	MG	DA	3117	1/1	0.94	0.34	5,5,5,5	0
57	MG	AA	1633	1/1	0.94	0.30	26,26,26,26	0
57	MG	CA	1705	1/1	0.94	0.11	5,5,5,5	0
57	MG	BA	3085	1/1	0.94	0.19	1,1,1,1	0
57	MG	BA	3444	1/1	0.94	0.65	15,15,15,15	1
57	MG	DA	3284	1/1	0.94	0.38	18,18,18,18	0
57	MG	DA	3205	1/1	0.94	0.08	13,13,13,13	0
57	MG	CA	1648	1/1	0.94	0.50	14,14,14,14	1
57	MG	CA	1755	1/1	0.94	0.30	1,1,1,1	0
57	MG	CA	1695	1/1	0.94	0.25	1,1,1,1	1
57	MG	DA	3067	1/1	0.95	0.32	1,1,1,1	0
57	MG	AA	1601	1/1	0.95	0.15	21,21,21,21	0
57	MG	DA	3418	1/1	0.95	0.62	24,24,24,24	1
57	MG	DA	3352	1/1	0.95	0.27	123,123,123,123	1
57	MG	BA	3113	1/1	0.95	0.29	1,1,1,1	0
57	MG	AA	1729	1/1	0.95	0.10	19,19,19,19	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3358	1/1	0.95	0.30	22,22,22,22	0
57	MG	B1	103	1/1	0.95	0.15	1,1,1,1	0
57	MG	BA	3037	1/1	0.95	0.25	27,27,27,27	0
57	MG	BA	3210	1/1	0.95	1.17	10,10,10,10	1
57	MG	BA	3235	1/1	0.95	0.44	22,22,22,22	0
57	MG	CA	1781	1/1	0.95	0.21	1,1,1,1	1
57	MG	CA	1624	1/1	0.95	0.69	40,40,40,40	0
57	MG	DA	3058	1/1	0.95	0.22	42,42,42,42	0
57	MG	DA	3410	1/1	0.95	0.16	31,31,31,31	1
57	MG	AA	1617	1/1	0.95	0.11	6,6,6,6	1
57	MG	DA	3148	1/1	0.95	0.21	26,26,26,26	0
57	MG	BA	3004	1/1	0.95	0.20	31,31,31,31	1
57	MG	CA	1647	1/1	0.95	0.42	82,82,82,82	0
57	MG	BA	3301	1/1	0.95	0.12	24,24,24,24	0
57	MG	DA	3280	1/1	0.95	0.16	24,24,24,24	1
57	MG	AA	1638	1/1	0.95	0.41	39,39,39,39	0
57	MG	DP	203	1/1	0.95	0.17	1,1,1,1	1
57	MG	DA	3001	1/1	0.95	0.23	49,49,49,49	0
57	MG	DA	3199	1/1	0.95	0.23	5,5,5,5	0
57	MG	DF	302	1/1	0.95	0.21	15,15,15,15	0
57	MG	BA	3329	1/1	0.95	0.56	61,61,61,61	0
57	MG	BA	3217	1/1	0.95	0.33	35,35,35,35	0
57	MG	DA	3059	1/1	0.95	0.18	1,1,1,1	0
57	MG	AA	1700	1/1	0.95	0.41	34,34,34,34	0
57	MG	DA	3260	1/1	0.95	0.31	8,8,8,8	1
57	MG	CA	1678	1/1	0.95	0.10	4,4,4,4	0
57	MG	DA	3411	1/1	0.95	0.17	19,19,19,19	0
57	MG	BA	3160	1/1	0.95	0.38	6,6,6,6	0
57	MG	DF	303	1/1	0.95	0.61	4,4,4,4	1
57	MG	AA	1681	1/1	0.95	0.31	14,14,14,14	0
57	MG	DA	3035	1/1	0.95	0.10	48,48,48,48	0
57	MG	DA	3437	1/1	0.95	0.27	59,59,59,59	1
57	MG	DA	3333	1/1	0.95	0.10	15,15,15,15	0
57	MG	DA	3356	1/1	0.95	0.17	57,57,57,57	0
57	MG	AA	1772	1/1	0.95	0.57	52,52,52,52	0
57	MG	AA	1614	1/1	0.95	0.59	11,11,11,11	0
57	MG	DA	3408	1/1	0.95	0.28	1,1,1,1	0
57	MG	BA	3079	1/1	0.95	0.20	1,1,1,1	0
57	MG	AX	105	1/1	0.95	0.14	35,35,35,35	0
57	MG	BA	3326	1/1	0.95	0.19	23,23,23,23	1
57	MG	BA	3008	1/1	0.95	0.34	19,19,19,19	0
57	MG	DA	3146	1/1	0.95	0.16	3,3,3,3	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DP	204	1/1	0.95	0.13	15,15,15,15	1
57	MG	DA	3040	1/1	0.95	0.13	1,1,1,1	0
57	MG	DA	3425	1/1	0.95	0.26	19,19,19,19	0
57	MG	DA	3213	1/1	0.95	0.24	24,24,24,24	0
57	MG	BX	101	1/1	0.95	0.21	26,26,26,26	1
57	MG	AW	101	1/1	0.95	1.05	51,51,51,51	1
57	MG	BA	3312	1/1	0.95	0.22	1,1,1,1	0
57	MG	BA	3126	1/1	0.95	0.33	1,1,1,1	0
57	MG	DA	3041	1/1	0.95	0.12	24,24,24,24	0
57	MG	BA	3044	1/1	0.95	0.38	36,36,36,36	0
57	MG	DA	3131	1/1	0.95	0.13	22,22,22,22	0
57	MG	DA	3007	1/1	0.95	0.16	14,14,14,14	0
57	MG	B1	101	1/1	0.95	0.30	27,27,27,27	0
57	MG	DA	3211	1/1	0.95	0.17	39,39,39,39	1
57	MG	DA	3009	1/1	0.95	0.27	1,1,1,1	1
57	MG	BA	3019	1/1	0.95	0.48	1,1,1,1	0
57	MG	DB	211	1/1	0.95	0.15	1,1,1,1	1
57	MG	DA	3216	1/1	0.95	0.10	28,28,28,28	1
57	MG	DA	3203	1/1	0.95	0.41	1,1,1,1	1
57	MG	AA	1796	1/1	0.95	0.19	9,9,9,9	0
57	MG	BA	3260	1/1	0.95	0.45	1,1,1,1	0
57	MG	CA	1634	1/1	0.95	0.42	25,25,25,25	0
57	MG	CA	1670	1/1	0.95	0.13	23,23,23,23	0
57	MG	DA	3032	1/1	0.95	0.19	23,23,23,23	0
57	MG	CA	1779	1/1	0.95	0.43	34,34,34,34	0
57	MG	DN	201	1/1	0.95	0.18	21,21,21,21	1
57	MG	B2	602	1/1	0.95	0.24	1,1,1,1	1
57	MG	CE	203	1/1	0.95	0.45	38,38,38,38	1
57	MG	CA	1671	1/1	0.95	0.33	21,21,21,21	0
57	MG	BA	3174	1/1	0.95	0.19	1,1,1,1	1
57	MG	AA	1642	1/1	0.95	0.14	26,26,26,26	0
57	MG	DA	3162	1/1	0.95	0.14	2,2,2,2	1
57	MG	AL	201	1/1	0.95	0.15	1,1,1,1	1
57	MG	BA	3199	1/1	0.95	0.46	5,5,5,5	1
57	MG	DA	3398	1/1	0.95	0.18	13,13,13,13	1
57	MG	CA	1669	1/1	0.95	0.41	32,32,32,32	0
57	MG	AA	1775	1/1	0.95	0.17	14,14,14,14	1
57	MG	BA	3192	1/1	0.95	0.44	15,15,15,15	0
57	MG	BA	3226	1/1	0.95	0.32	1,1,1,1	1
57	MG	BA	3426	1/1	0.95	0.20	27,27,27,27	0
57	MG	DA	3391	1/1	0.95	0.12	19,19,19,19	0
57	MG	AA	1709	1/1	0.95	0.18	3,3,3,3	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3015	1/1	0.95	0.15	6,6,6,6	0
57	MG	DA	3394	1/1	0.95	0.17	30,30,30,30	1
57	MG	BA	3121	1/1	0.95	0.35	30,30,30,30	0
57	MG	DA	3013	1/1	0.95	0.51	7,7,7,7	1
57	MG	CA	1635	1/1	0.95	0.06	6,6,6,6	0
57	MG	BA	3288	1/1	0.95	0.22	44,44,44,44	1
57	MG	DA	3068	1/1	0.95	0.26	24,24,24,24	0
57	MG	DB	214	1/1	0.96	0.95	25,25,25,25	1
57	MG	BA	3432	1/1	0.96	0.33	47,47,47,47	0
57	MG	BA	3378	1/1	0.96	0.15	31,31,31,31	0
57	MG	BA	3066	1/1	0.96	0.14	20,20,20,20	0
57	MG	BA	3268	1/1	0.96	0.44	30,30,30,30	0
57	MG	BA	3202	1/1	0.96	0.21	14,14,14,14	0
57	MG	DA	3250	1/1	0.96	0.18	1,1,1,1	0
57	MG	AA	1612	1/1	0.96	0.26	36,36,36,36	0
57	MG	CA	1606	1/1	0.96	0.43	29,29,29,29	0
57	MG	AA	1771	1/1	0.96	0.17	9,9,9,9	0
57	MG	CA	1750	1/1	0.96	0.36	14,14,14,14	0
57	MG	BA	3183	1/1	0.96	0.76	52,52,52,52	0
57	MG	AA	1607	1/1	0.96	0.16	1,1,1,1	0
57	MG	BA	3063	1/1	0.96	0.13	6,6,6,6	0
57	MG	CA	1604	1/1	0.96	0.10	41,41,41,41	0
57	MG	AA	1708	1/1	0.96	0.19	37,37,37,37	0
57	MG	DA	3168	1/1	0.96	0.87	36,36,36,36	1
57	MG	DA	3379	1/1	0.96	0.16	47,47,47,47	0
57	MG	BA	3255	1/1	0.96	0.17	25,25,25,25	0
57	MG	CA	1628	1/1	0.96	0.68	35,35,35,35	0
57	MG	BA	3196	1/1	0.96	0.44	19,19,19,19	0
57	MG	DA	3133	1/1	0.96	0.16	9,9,9,9	0
57	MG	BE	301	1/1	0.96	0.22	1,1,1,1	1
57	MG	AA	1731	1/1	0.96	0.12	79,79,79,79	1
57	MG	BA	3430	1/1	0.96	0.29	1,1,1,1	1
57	MG	CA	1751	1/1	0.96	0.31	10,10,10,10	0
57	MG	DA	3364	1/1	0.96	0.27	42,42,42,42	0
57	MG	DU	202	1/1	0.96	0.68	1,1,1,1	1
57	MG	BF	301	1/1	0.96	0.38	14,14,14,14	0
57	MG	CA	1637	1/1	0.96	0.15	47,47,47,47	0
57	MG	CA	1611	1/1	0.96	0.10	22,22,22,22	0
57	MG	DA	3124	1/1	0.96	0.22	24,24,24,24	0
57	MG	BA	3293	1/1	0.96	0.10	19,19,19,19	0
57	MG	DA	3294	1/1	0.96	0.18	73,73,73,73	1
57	MG	AA	1635	1/1	0.96	0.24	2,2,2,2	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3237	1/1	0.96	0.34	22,22,22,22	0
57	MG	BA	3363	1/1	0.96	0.12	90,90,90,90	0
57	MG	BA	3305	1/1	0.96	0.09	37,37,37,37	0
57	MG	DA	3266	1/1	0.96	0.20	17,17,17,17	1
57	MG	AM	201	1/1	0.96	0.12	28,28,28,28	0
57	MG	BA	3400	1/1	0.96	0.22	1,1,1,1	0
57	MG	BA	3181	1/1	0.96	0.08	17,17,17,17	0
57	MG	CA	1630	1/1	0.96	0.33	40,40,40,40	0
57	MG	AA	1677	1/1	0.96	0.12	42,42,42,42	0
57	MG	DA	3374	1/1	0.96	0.15	28,28,28,28	0
57	MG	DA	3034	1/1	0.96	0.29	33,33,33,33	0
57	MG	BA	3398	1/1	0.96	0.10	11,11,11,11	0
57	MG	BA	3111	1/1	0.96	0.06	1,1,1,1	0
57	MG	BA	3133	1/1	0.96	0.22	1,1,1,1	0
57	MG	AG	201	1/1	0.96	0.22	5,5,5,5	1
57	MG	DA	3134	1/1	0.96	0.31	8,8,8,8	0
57	MG	BA	3114	1/1	0.96	0.27	1,1,1,1	0
57	MG	CA	1625	1/1	0.96	0.17	23,23,23,23	0
57	MG	AA	1626	1/1	0.96	0.09	1,1,1,1	0
57	MG	BA	3190	1/1	0.96	0.14	17,17,17,17	0
57	MG	BA	3155	1/1	0.96	0.16	37,37,37,37	0
57	MG	DA	3024	1/1	0.96	0.32	32,32,32,32	0
57	MG	DA	3384	1/1	0.96	0.16	1,1,1,1	1
57	MG	AA	1627	1/1	0.96	0.48	9,9,9,9	0
57	MG	CA	1682	1/1	0.96	0.24	9,9,9,9	0
57	MG	BA	3171	1/1	0.96	0.17	1,1,1,1	1
57	MG	DA	3094	1/1	0.96	0.59	1,1,1,1	1
57	MG	DA	3288	1/1	0.96	0.26	7,7,7,7	1
57	MG	BA	3361	1/1	0.96	0.29	1,1,1,1	1
57	MG	BA	3302	1/1	0.96	0.05	19,19,19,19	1
57	MG	CA	1642	1/1	0.96	0.13	2,2,2,2	0
57	MG	BA	3427	1/1	0.96	0.29	99,99,99,99	1
57	MG	CA	1685	1/1	0.96	0.22	25,25,25,25	1
57	MG	AA	1624	1/1	0.96	0.24	9,9,9,9	0
57	MG	CA	1752	1/1	0.96	0.23	85,85,85,85	0
57	MG	BB	212	1/1	0.96	0.15	1,1,1,1	1
57	MG	DA	3196	1/1	0.96	0.35	7,7,7,7	1
57	MG	CA	1734	1/1	0.96	0.21	46,46,46,46	1
57	MG	DA	3081	1/1	0.96	0.14	11,11,11,11	0
57	MG	CA	1701	1/1	0.96	0.24	3,3,3,3	0
57	MG	DA	3300	1/1	0.96	0.19	38,38,38,38	0
57	MG	DA	3036	1/1	0.96	0.46	13,13,13,13	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3025	1/1	0.96	0.31	3,3,3,3	0
57	MG	CL	201	1/1	0.96	0.09	1,1,1,1	0
57	MG	DA	3155	1/1	0.96	0.37	21,21,21,21	0
57	MG	AA	1685	1/1	0.96	0.13	12,12,12,12	0
57	MG	BA	3300	1/1	0.96	0.28	1,1,1,1	1
57	MG	DA	3118	1/1	0.96	0.20	9,9,9,9	0
57	MG	CV	103	1/1	0.96	0.05	44,44,44,44	0
57	MG	DA	3003	1/1	0.96	0.09	1,1,1,1	1
57	MG	DA	3051	1/1	0.96	0.18	1,1,1,1	0
57	MG	DA	3283	1/1	0.96	0.21	7,7,7,7	1
57	MG	DA	3171	1/1	0.96	0.18	33,33,33,33	0
57	MG	AL	202	1/1	0.96	0.19	1,1,1,1	0
57	MG	AA	1756	1/1	0.96	0.15	58,58,58,58	0
57	MG	DA	3404	1/1	0.96	0.25	43,43,43,43	0
57	MG	AA	1804	1/1	0.96	0.26	20,20,20,20	0
57	MG	AA	1619	1/1	0.96	0.16	25,25,25,25	0
57	MG	DA	3433	1/1	0.96	0.18	28,28,28,28	1
57	MG	DA	3113	1/1	0.96	0.14	1,1,1,1	0
57	MG	DB	217	1/1	0.96	0.05	53,53,53,53	1
57	MG	AW	105	1/1	0.96	0.51	82,82,82,82	0
57	MG	BA	3028	1/1	0.96	0.41	1,1,1,1	0
57	MG	DF	301	1/1	0.96	0.19	38,38,38,38	1
57	MG	DA	3232	1/1	0.96	0.42	1,1,1,1	0
57	MG	BA	3146	1/1	0.96	0.13	1,1,1,1	0
57	MG	DX	102	1/1	0.96	0.26	23,23,23,23	1
57	MG	DA	3249	1/1	0.96	0.11	1,1,1,1	0
57	MG	CA	1740	1/1	0.96	0.35	44,44,44,44	0
57	MG	DA	3191	1/1	0.96	0.19	5,5,5,5	1
57	MG	CA	1638	1/1	0.96	0.11	6,6,6,6	0
57	MG	DA	3065	1/1	0.96	0.26	1,1,1,1	1
57	MG	DA	3306	1/1	0.96	0.14	1,1,1,1	0
57	MG	BA	3377	1/1	0.96	0.26	95,95,95,95	0
57	MG	AA	1776	1/1	0.96	0.20	33,33,33,33	1
57	MG	AA	1690	1/1	0.96	0.15	40,40,40,40	0
57	MG	BA	3033	1/1	0.96	0.31	20,20,20,20	0
57	MG	CA	1790	1/1	0.96	0.18	19,19,19,19	0
57	MG	DV	201	1/1	0.96	0.15	1,1,1,1	0
57	MG	DA	3293	1/1	0.96	0.32	1,1,1,1	1
57	MG	DA	3399	1/1	0.96	0.14	19,19,19,19	0
57	MG	BP	202	1/1	0.96	0.54	9,9,9,9	1
57	MG	BA	3233	1/1	0.96	0.30	1,1,1,1	0
57	MG	BA	3149	1/1	0.96	0.18	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BB	207	1/1	0.96	0.04	11,11,11,11	0
57	MG	DA	3298	1/1	0.96	0.08	22,22,22,22	0
57	MG	BA	3177	1/1	0.96	0.26	36,36,36,36	1
57	MG	AA	1673	1/1	0.96	0.12	12,12,12,12	1
57	MG	BA	3070	1/1	0.96	0.27	24,24,24,24	0
57	MG	BA	3110	1/1	0.96	0.19	32,32,32,32	1
57	MG	AA	1774	1/1	0.96	0.30	23,23,23,23	0
57	MG	BA	3003	1/1	0.96	0.34	68,68,68,68	1
57	MG	BP	201	1/1	0.96	0.28	11,11,11,11	1
57	MG	CA	1636	1/1	0.96	0.23	1,1,1,1	0
57	MG	DA	3287	1/1	0.96	0.14	28,28,28,28	0
57	MG	DA	3017	1/1	0.96	0.07	11,11,11,11	0
57	MG	DA	3055	1/1	0.96	0.08	1,1,1,1	0
57	MG	CA	1633	1/1	0.97	0.07	2,2,2,2	0
57	MG	BA	3065	1/1	0.97	0.18	7,7,7,7	0
57	MG	DA	3401	1/1	0.97	0.29	49,49,49,49	1
57	MG	DA	3114	1/1	0.97	0.13	1,1,1,1	0
57	MG	BA	3108	1/1	0.97	0.09	25,25,25,25	0
57	MG	DA	3056	1/1	0.97	0.24	11,11,11,11	0
57	MG	DA	3031	1/1	0.97	0.28	2,2,2,2	0
57	MG	DU	201	1/1	0.97	0.29	77,77,77,77	1
57	MG	CA	1699	1/1	0.97	0.50	6,6,6,6	1
57	MG	DA	3386	1/1	0.97	0.14	19,19,19,19	0
57	MG	CA	1791	1/1	0.97	0.13	1,1,1,1	1
57	MG	AA	1801	1/1	0.97	0.35	1,1,1,1	0
57	MG	BA	3072	1/1	0.97	0.34	100,100,100,100	0
57	MG	BA	3438	1/1	0.97	0.09	44,44,44,44	0
57	MG	DA	3110	1/1	0.97	0.15	1,1,1,1	1
57	MG	CA	1753	1/1	0.97	0.20	7,7,7,7	1
57	MG	BA	3061	1/1	0.97	0.21	3,3,3,3	0
57	MG	AA	1658	1/1	0.97	0.24	1,1,1,1	0
57	MG	DA	3128	1/1	0.97	0.12	1,1,1,1	1
57	MG	BA	3022	1/1	0.97	0.47	1,1,1,1	0
57	MG	BA	3410	1/1	0.97	0.20	6,6,6,6	1
57	MG	DA	3158	1/1	0.97	0.13	23,23,23,23	0
57	MG	CX	102	1/1	0.97	0.17	1,1,1,1	1
57	MG	DA	3186	1/1	0.97	0.17	24,24,24,24	0
57	MG	B2	601	1/1	0.97	0.51	22,22,22,22	0
57	MG	DA	3427	1/1	0.97	0.22	1,1,1,1	0
57	MG	DA	3190	1/1	0.97	0.41	32,32,32,32	0
57	MG	BA	3172	1/1	0.97	0.29	100,100,100,100	1
57	MG	BA	3012	1/1	0.97	0.12	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1606	1/1	0.97	0.15	1,1,1,1	0
57	MG	AA	1768	1/1	0.97	0.23	7,7,7,7	0
57	MG	AA	1722	1/1	0.97	0.12	37,37,37,37	0
57	MG	B5	101	1/1	0.97	0.23	1,1,1,1	0
57	MG	DA	3043	1/1	0.97	0.17	6,6,6,6	0
57	MG	AA	1664	1/1	0.97	0.21	4,4,4,4	0
57	MG	BA	3127	1/1	0.97	0.44	3,3,3,3	1
57	MG	DA	3129	1/1	0.97	0.18	1,1,1,1	0
57	MG	BA	3194	1/1	0.97	0.42	19,19,19,19	0
57	MG	BA	3173	1/1	0.97	0.29	24,24,24,24	0
57	MG	BA	3090	1/1	0.97	0.25	24,24,24,24	1
57	MG	AA	1795	1/1	0.97	0.16	1,1,1,1	0
57	MG	BA	3314	1/1	0.97	0.48	23,23,23,23	0
57	MG	BA	3136	1/1	0.97	0.29	1,1,1,1	0
57	MG	BD	301	1/1	0.97	0.27	1,1,1,1	0
57	MG	BA	3122	1/1	0.97	0.20	1,1,1,1	1
57	MG	DA	3052	1/1	0.97	0.16	1,1,1,1	0
57	MG	D1	101	1/1	0.97	0.46	55,55,55,55	0
57	MG	CA	1688	1/1	0.97	0.37	21,21,21,21	0
57	MG	CA	1714	1/1	0.97	0.10	53,53,53,53	0
57	MG	BA	3296	1/1	0.97	0.26	11,11,11,11	0
57	MG	BA	3098	1/1	0.97	0.15	1,1,1,1	0
57	MG	BA	3104	1/1	0.97	0.10	1,1,1,1	0
57	MG	AX	102	1/1	0.97	0.19	4,4,4,4	0
57	MG	DA	3029	1/1	0.97	0.14	1,1,1,1	0
57	MG	BA	3119	1/1	0.97	0.18	4,4,4,4	0
57	MG	BA	3049	1/1	0.97	0.20	17,17,17,17	0
57	MG	BU	202	1/1	0.97	0.25	18,18,18,18	1
57	MG	DA	3025	1/1	0.97	0.19	5,5,5,5	0
57	MG	DA	3027	1/1	0.97	0.48	55,55,55,55	1
57	MG	B1	104	1/1	0.97	0.18	62,62,62,62	0
57	MG	DA	3130	1/1	0.97	0.09	33,33,33,33	0
57	MG	BA	3382	1/1	0.97	0.06	38,38,38,38	0
57	MG	DA	3093	1/1	0.97	0.10	9,9,9,9	0
57	MG	BA	3352	1/1	0.97	0.94	44,44,44,44	1
57	MG	DA	3151	1/1	0.97	0.20	1,1,1,1	1
57	MG	BA	3242	1/1	0.97	0.11	1,1,1,1	0
57	MG	DA	3167	1/1	0.97	0.10	12,12,12,12	0
57	MG	DA	3304	1/1	0.97	0.46	1,1,1,1	1
57	MG	DA	3308	1/1	0.97	0.58	3,3,3,3	0
57	MG	AA	1784	1/1	0.97	0.42	43,43,43,43	1
57	MG	BA	3185	1/1	0.97	0.09	1,1,1,1	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3088	1/1	0.97	0.51	1,1,1,1	1
57	MG	DA	3193	1/1	0.97	0.12	1,1,1,1	1
57	MG	DA	3311	1/1	0.97	0.22	11,11,11,11	1
57	MG	DA	3413	1/1	0.97	0.10	10,10,10,10	1
57	MG	AA	1628	1/1	0.97	0.14	1,1,1,1	1
57	MG	CA	1777	1/1	0.97	0.12	47,47,47,47	0
57	MG	BA	3354	1/1	0.97	0.14	58,58,58,58	0
57	MG	D5	101	1/1	0.97	0.11	20,20,20,20	0
57	MG	AA	1745	1/1	0.97	0.35	79,79,79,79	0
57	MG	DA	3053	1/1	0.97	0.28	3,3,3,3	0
57	MG	BA	3051	1/1	0.97	0.56	1,1,1,1	0
57	MG	AA	1699	1/1	0.97	0.07	1,1,1,1	1
57	MG	AA	1786	1/1	0.97	0.12	1,1,1,1	1
57	MG	BA	3076	1/1	0.97	0.20	1,1,1,1	0
57	MG	AA	1636	1/1	0.97	0.06	16,16,16,16	0
57	MG	AW	107	1/1	0.97	0.77	57,57,57,57	0
57	MG	DA	3209	1/1	0.97	0.24	1,1,1,1	1
57	MG	DA	3429	1/1	0.97	0.76	36,36,36,36	1
57	MG	BA	3304	1/1	0.97	0.14	24,24,24,24	0
57	MG	BA	3208	1/1	0.97	0.21	23,23,23,23	1
57	MG	DA	3023	1/1	0.97	0.41	1,1,1,1	0
57	MG	BA	3428	1/1	0.97	0.11	1,1,1,1	0
57	MG	BB	211	1/1	0.97	0.31	44,44,44,44	1
57	MG	BD	303	1/1	0.97	0.39	13,13,13,13	1
57	MG	DB	203	1/1	0.97	0.29	60,60,60,60	0
57	MG	DA	3063	1/1	0.97	0.44	1,1,1,1	0
57	MG	AA	1778	1/1	0.97	0.09	16,16,16,16	1
57	MG	BA	3082	1/1	0.97	0.17	1,1,1,1	1
57	MG	BA	3342	1/1	0.97	0.16	26,26,26,26	1
57	MG	DA	3112	1/1	0.97	0.41	1,1,1,1	0
57	MG	AA	1637	1/1	0.97	0.22	11,11,11,11	0
57	MG	CA	1708	1/1	0.97	0.15	99,99,99,99	0
57	MG	BA	3047	1/1	0.97	0.25	1,1,1,1	0
57	MG	DA	3297	1/1	0.97	0.35	11,11,11,11	1
57	MG	DA	3327	1/1	0.97	0.07	2,2,2,2	1
57	MG	DA	3240	1/1	0.97	0.13	72,72,72,72	0
57	MG	DA	3383	1/1	0.97	0.15	42,42,42,42	1
57	MG	CA	1662	1/1	0.97	0.44	1,1,1,1	0
57	MG	BA	3313	1/1	0.97	0.40	18,18,18,18	0
57	MG	DA	3313	1/1	0.97	0.28	5,5,5,5	0
57	MG	BA	3036	1/1	0.97	0.40	18,18,18,18	0
57	MG	DA	3261	1/1	0.97	0.37	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3403	1/1	0.97	0.22	55,55,55,55	1
57	MG	BA	3186	1/1	0.97	0.17	5,5,5,5	1
57	MG	DA	3070	1/1	0.97	0.35	40,40,40,40	1
57	MG	CA	1749	1/1	0.97	0.10	49,49,49,49	1
57	MG	CA	1691	1/1	0.97	0.09	1,1,1,1	0
57	MG	DA	3393	1/1	0.97	0.14	1,1,1,1	1
57	MG	B7	101	1/1	0.97	0.08	8,8,8,8	0
57	MG	DA	3039	1/1	0.97	0.14	2,2,2,2	0
57	MG	BA	3256	1/1	0.97	0.18	1,1,1,1	1
57	MG	CA	1676	1/1	0.97	0.21	43,43,43,43	1
57	MG	DA	3097	1/1	0.97	0.12	55,55,55,55	1
57	MG	BA	3193	1/1	0.97	0.14	1,1,1,1	0
57	MG	DA	3200	1/1	0.97	0.17	2,2,2,2	0
57	MG	BA	3054	1/1	0.97	0.14	20,20,20,20	0
57	MG	DA	3291	1/1	0.97	0.16	1,1,1,1	1
57	MG	BA	3075	1/1	0.97	0.35	6,6,6,6	1
57	MG	BA	3131	1/1	0.97	0.12	1,1,1,1	0
57	MG	BA	3039	1/1	0.97	0.09	1,1,1,1	0
57	MG	AA	1672	1/1	0.97	0.17	24,24,24,24	0
57	MG	BA	3433	1/1	0.97	0.14	33,33,33,33	0
57	MG	BA	3059	1/1	0.97	0.12	9,9,9,9	0
57	MG	BA	3023	1/1	0.97	0.10	1,1,1,1	1
57	MG	AW	104	1/1	0.97	0.09	14,14,14,14	1
57	MG	CA	1766	1/1	0.97	0.16	1,1,1,1	0
57	MG	CA	1787	1/1	0.97	0.48	11,11,11,11	0
57	MG	DA	3157	1/1	0.97	0.25	37,37,37,37	0
57	MG	BA	3115	1/1	0.97	0.11	3,3,3,3	0
57	MG	BA	3148	1/1	0.97	0.28	8,8,8,8	0
57	MG	DA	3314	1/1	0.97	0.37	1,1,1,1	1
57	MG	D7	101	1/1	0.97	0.12	93,93,93,93	1
57	MG	BA	3221	1/1	0.97	0.16	49,49,49,49	0
57	MG	BA	3106	1/1	0.97	0.17	19,19,19,19	1
57	MG	AA	1618	1/1	0.97	0.17	1,1,1,1	0
57	MG	AA	1750	1/1	0.97	0.12	42,42,42,42	0
57	MG	CA	1614	1/1	0.97	0.33	6,6,6,6	0
57	MG	CA	1621	1/1	0.97	0.16	31,31,31,31	0
57	MG	DA	3092	1/1	0.97	0.41	1,1,1,1	0
57	MG	BA	3046	1/1	0.97	0.12	15,15,15,15	0
57	MG	BA	3097	1/1	0.98	0.10	1,1,1,1	0
57	MG	DA	3412	1/1	0.98	0.29	12,12,12,12	0
57	MG	AA	1670	1/1	0.98	0.20	22,22,22,22	0
57	MG	DA	3021	1/1	0.98	0.19	2,2,2,2	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3105	1/1	0.98	0.17	1,1,1,1	0
57	MG	BA	3317	1/1	0.98	0.12	60,60,60,60	0
57	MG	DA	3230	1/1	0.98	0.13	35,35,35,35	0
57	MG	BA	3032	1/1	0.98	0.15	1,1,1,1	0
57	MG	DA	3098	1/1	0.98	0.13	5,5,5,5	0
57	MG	BA	3419	1/1	0.98	0.07	57,57,57,57	1
57	MG	DA	3305	1/1	0.98	0.21	1,1,1,1	0
57	MG	CA	1773	1/1	0.98	0.30	9,9,9,9	0
57	MG	DA	3099	1/1	0.98	0.29	1,1,1,1	1
57	MG	DA	3188	1/1	0.98	0.19	26,26,26,26	0
57	MG	AA	1697	1/1	0.98	0.21	17,17,17,17	1
57	MG	BA	3138	1/1	0.98	0.21	12,12,12,12	0
57	MG	BA	3116	1/1	0.98	0.13	13,13,13,13	0
57	MG	BA	3200	1/1	0.98	0.14	5,5,5,5	1
57	MG	AA	1660	1/1	0.98	0.32	17,17,17,17	0
57	MG	DA	3256	1/1	0.98	0.38	1,1,1,1	0
57	MG	CA	1608	1/1	0.98	0.11	17,17,17,17	0
57	MG	DA	3096	1/1	0.98	0.29	1,1,1,1	0
57	MG	DA	3315	1/1	0.98	0.32	6,6,6,6	0
57	MG	BA	3042	1/1	0.98	0.18	1,1,1,1	0
57	MG	AA	1751	1/1	0.98	0.23	14,14,14,14	0
57	MG	DA	3108	1/1	0.98	0.19	13,13,13,13	0
57	MG	DA	3353	1/1	0.98	0.25	107,107,107,107	1
57	MG	CA	1789	1/1	0.98	0.25	73,73,73,73	0
57	MG	DA	3076	1/1	0.98	0.33	1,1,1,1	0
57	MG	AA	1647	1/1	0.98	0.10	17,17,17,17	1
57	MG	AA	1632	1/1	0.98	0.19	4,4,4,4	0
57	MG	DA	3380	1/1	0.98	0.10	16,16,16,16	1
57	MG	CA	1723	1/1	0.98	0.19	10,10,10,10	1
57	MG	CA	1619	1/1	0.98	0.07	3,3,3,3	0
57	MG	CA	1626	1/1	0.98	0.09	23,23,23,23	0
57	MG	DA	3272	1/1	0.98	0.17	5,5,5,5	0
57	MG	DA	3178	1/1	0.98	0.15	1,1,1,1	1
57	MG	AA	1654	1/1	0.98	0.36	65,65,65,65	0
57	MG	DA	3062	1/1	0.98	0.54	21,21,21,21	0
57	MG	DA	3198	1/1	0.98	0.16	11,11,11,11	0
57	MG	AA	1727	1/1	0.98	0.14	63,63,63,63	0
57	MG	CA	1732	1/1	0.98	0.27	10,10,10,10	0
57	MG	AA	1692	1/1	0.98	0.19	3,3,3,3	1
57	MG	BA	3092	1/1	0.98	0.17	5,5,5,5	0
57	MG	DA	3254	1/1	0.98	0.24	109,109,109,109	1
57	MG	BA	3223	1/1	0.98	0.22	9,9,9,9	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3244	1/1	0.98	0.13	57,57,57,57	0
57	MG	DA	3348	1/1	0.98	0.12	9,9,9,9	1
57	MG	DA	3169	1/1	0.98	0.34	1,1,1,1	1
57	MG	AA	1678	1/1	0.98	0.07	13,13,13,13	0
57	MG	BA	3084	1/1	0.98	0.17	1,1,1,1	1
57	MG	DA	3153	1/1	0.98	0.15	1,1,1,1	0
57	MG	CA	1667	1/1	0.98	0.26	8,8,8,8	0
57	MG	BA	3384	1/1	0.98	0.15	1,1,1,1	1
57	MG	CA	1665	1/1	0.98	0.48	34,34,34,34	0
57	MG	BA	3249	1/1	0.98	0.40	58,58,58,58	0
57	MG	BA	3265	1/1	0.98	0.11	37,37,37,37	0
57	MG	AA	1714	1/1	0.98	0.18	23,23,23,23	0
57	MG	DA	3142	1/1	0.98	0.28	21,21,21,21	0
57	MG	CA	1710	1/1	0.98	0.18	16,16,16,16	0
57	MG	BA	3278	1/1	0.98	0.07	8,8,8,8	0
57	MG	DA	3077	1/1	0.98	0.11	1,1,1,1	0
57	MG	DA	3074	1/1	0.98	0.07	2,2,2,2	0
57	MG	BD	302	1/1	0.98	0.26	1,1,1,1	0
57	MG	DA	3105	1/1	0.98	0.10	26,26,26,26	0
57	MG	DA	3109	1/1	0.98	0.08	17,17,17,17	0
57	MG	AA	1657	1/1	0.98	0.16	8,8,8,8	1
57	MG	BA	3151	1/1	0.98	0.26	1,1,1,1	0
57	MG	BA	3240	1/1	0.98	0.20	8,8,8,8	0
57	MG	DA	3416	1/1	0.98	0.14	1,1,1,1	1
57	MG	BA	3261	1/1	0.98	0.43	1,1,1,1	0
57	MG	CA	1799	1/1	0.98	0.29	39,39,39,39	1
57	MG	BA	3344	1/1	0.98	0.17	1,1,1,1	1
57	MG	D2	2602	1/1	0.98	0.34	6,6,6,6	1
57	MG	DA	3397	1/1	0.98	0.31	1,1,1,1	0
57	MG	DF	304	1/1	0.98	0.23	66,66,66,66	0
57	MG	BA	3071	1/1	0.98	0.14	20,20,20,20	0
57	MG	BA	3414	1/1	0.98	0.17	20,20,20,20	0
57	MG	BA	3254	1/1	0.98	0.06	5,5,5,5	0
57	MG	BN	201	1/1	0.98	0.19	47,47,47,47	1
57	MG	BP	203	1/1	0.98	0.31	1,1,1,1	1
58	ZN	CN	102	1/1	0.98	0.14	114,114,114,114	0
57	MG	CA	1663	1/1	0.98	0.56	15,15,15,15	1
57	MG	BA	3087	1/1	0.98	0.09	18,18,18,18	0
57	MG	DX	101	1/1	0.98	0.10	52,52,52,52	1
57	MG	DA	3104	1/1	0.98	0.14	9,9,9,9	1
57	MG	DA	3184	1/1	0.98	0.19	1,1,1,1	0
57	MG	BA	3034	1/1	0.98	0.14	15,15,15,15	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	DA	3091	1/1	0.98	0.14	1,1,1,1	0
57	MG	BA	3027	1/1	0.98	0.26	1,1,1,1	1
57	MG	DA	3143	1/1	0.98	0.39	24,24,24,24	0
57	MG	DA	3350	1/1	0.98	0.55	73,73,73,73	1
57	MG	BA	3021	1/1	0.98	0.16	2,2,2,2	0
57	MG	DA	3066	1/1	0.98	0.31	1,1,1,1	0
57	MG	DA	3111	1/1	0.98	0.14	4,4,4,4	0
57	MG	CA	1689	1/1	0.98	0.09	87,87,87,87	0
57	MG	DA	3197	1/1	0.98	0.24	3,3,3,3	0
57	MG	DA	3255	1/1	0.98	0.07	1,1,1,1	0
57	MG	DA	3119	1/1	0.98	0.31	1,1,1,1	0
57	MG	BA	3083	1/1	0.98	0.15	18,18,18,18	0
57	MG	DA	3334	1/1	0.98	0.21	30,30,30,30	0
57	MG	BA	3187	1/1	0.98	0.09	1,1,1,1	0
57	MG	BA	3311	1/1	0.98	0.12	1,1,1,1	0
57	MG	AA	1605	1/1	0.98	0.43	3,3,3,3	0
57	MG	B3	101	1/1	0.98	0.20	14,14,14,14	0
57	MG	AX	103	1/1	0.98	0.09	1,1,1,1	1
57	MG	DA	3132	1/1	0.98	0.08	34,34,34,34	0
57	MG	BA	3259	1/1	0.98	0.08	1,1,1,1	1
57	MG	DA	3137	1/1	0.98	0.15	5,5,5,5	1
57	MG	CA	1722	1/1	0.98	0.09	1,1,1,1	1
57	MG	DA	3204	1/1	0.98	0.21	51,51,51,51	1
57	MG	DA	3106	1/1	0.98	0.36	1,1,1,1	0
57	MG	DA	3375	1/1	0.98	0.52	74,74,74,74	1
57	MG	DA	3229	1/1	0.98	0.24	1,1,1,1	0
57	MG	BA	3142	1/1	0.98	0.15	67,67,67,67	0
57	MG	CX	104	1/1	0.98	0.14	39,39,39,39	0
57	MG	DA	3047	1/1	0.98	0.14	1,1,1,1	0
57	MG	BA	3048	1/1	0.98	0.46	1,1,1,1	0
57	MG	DA	3085	1/1	0.98	0.17	1,1,1,1	0
57	MG	AA	1641	1/1	0.98	0.14	26,26,26,26	0
57	MG	DA	3054	1/1	0.98	0.10	1,1,1,1	0
57	MG	BA	3064	1/1	0.98	0.30	1,1,1,1	0
57	MG	DA	3351	1/1	0.98	0.28	11,11,11,11	1
57	MG	BA	3080	1/1	0.98	0.23	26,26,26,26	0
57	MG	BA	3357	1/1	0.98	0.17	105,105,105,105	1
57	MG	BA	3211	1/1	0.98	0.17	50,50,50,50	1
57	MG	BA	3055	1/1	0.98	0.16	1,1,1,1	0
57	MG	BA	3355	1/1	0.98	0.10	37,37,37,37	0
57	MG	DA	3075	1/1	0.98	0.15	1,1,1,1	0
57	MG	DA	3050	1/1	0.98	0.44	136,136,136,136	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3045	1/1	0.98	0.24	1,1,1,1	1
57	MG	CA	1661	1/1	0.98	0.43	17,17,17,17	0
57	MG	DA	3335	1/1	0.98	0.12	7,7,7,7	1
57	MG	DA	3160	1/1	0.98	0.18	1,1,1,1	1
57	MG	AA	1662	1/1	0.98	0.09	41,41,41,41	1
57	MG	AA	1661	1/1	0.98	0.09	10,10,10,10	1
57	MG	BA	3147	1/1	0.98	0.72	1,1,1,1	1
57	MG	BA	3237	1/1	0.98	0.13	8,8,8,8	1
57	MG	BA	3424	1/1	0.98	0.15	37,37,37,37	0
57	MG	DA	3060	1/1	0.98	0.23	1,1,1,1	0
57	MG	DA	3082	1/1	0.98	0.14	20,20,20,20	0
57	MG	BA	3165	1/1	0.98	0.09	4,4,4,4	1
57	MG	CA	1627	1/1	0.98	0.09	5,5,5,5	0
57	MG	DA	3089	1/1	0.98	0.21	3,3,3,3	0
57	MG	DA	3246	1/1	0.98	0.23	57,57,57,57	0
57	MG	BA	3371	1/1	0.98	0.28	53,53,53,53	1
57	MG	BA	3041	1/1	0.98	0.18	46,46,46,46	1
57	MG	BA	3024	1/1	0.98	0.46	44,44,44,44	1
57	MG	BA	3077	1/1	0.98	0.14	1,1,1,1	0
57	MG	CA	1603	1/1	0.98	0.21	77,77,77,77	0
57	MG	BA	3234	1/1	0.98	0.18	12,12,12,12	0
57	MG	DA	3173	1/1	0.98	0.35	52,52,52,52	1
57	MG	DA	3251	1/1	0.99	0.23	10,10,10,10	1
57	MG	DA	3175	1/1	0.99	0.26	43,43,43,43	1
57	MG	BA	3218	1/1	0.99	0.08	50,50,50,50	0
57	MG	DD	302	1/1	0.99	0.13	2,2,2,2	0
57	MG	DA	3222	1/1	0.99	0.23	3,3,3,3	1
57	MG	BA	3067	1/1	0.99	0.12	1,1,1,1	1
57	MG	CA	1807	1/1	0.99	0.21	1,1,1,1	0
57	MG	DA	3183	1/1	0.99	0.11	11,11,11,11	0
58	ZN	AN	101	1/1	0.99	0.23	143,143,143,143	0
57	MG	BA	3243	1/1	0.99	0.25	1,1,1,1	0
57	MG	DA	3381	1/1	0.99	0.16	65,65,65,65	0
57	MG	DA	3421	1/1	0.99	0.44	40,40,40,40	0
57	MG	DA	3002	1/1	0.99	0.12	6,6,6,6	1
57	MG	AA	1630	1/1	0.99	0.16	19,19,19,19	0
57	MG	BA	3081	1/1	0.99	0.14	1,1,1,1	0
57	MG	AA	1675	1/1	0.99	0.29	41,41,41,41	1
57	MG	BA	3062	1/1	0.99	0.24	1,1,1,1	0
57	MG	AX	104	1/1	0.99	0.11	55,55,55,55	1
57	MG	DA	3116	1/1	0.99	0.40	1,1,1,1	1
57	MG	BA	3195	1/1	0.99	0.13	39,39,39,39	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	AA	1770	1/1	0.99	0.10	16,16,16,16	1
57	MG	DA	3015	1/1	0.99	0.05	1,1,1,1	1
57	MG	CX	101	1/1	0.99	0.08	56,56,56,56	1
57	MG	BA	3420	1/1	0.99	0.12	10,10,10,10	0
57	MG	BB	201	1/1	0.99	0.06	32,32,32,32	1
57	MG	BA	3273	1/1	0.99	0.14	57,57,57,57	1
57	MG	BA	3056	1/1	0.99	0.09	1,1,1,1	0
57	MG	DA	3263	1/1	0.99	0.15	16,16,16,16	0
57	MG	BA	3236	1/1	0.99	0.14	1,1,1,1	0
57	MG	DA	3061	1/1	0.99	0.08	31,31,31,31	0
57	MG	BA	3376	1/1	0.99	0.05	15,15,15,15	0
57	MG	DA	3243	1/1	0.99	0.09	18,18,18,18	0
57	MG	DA	3095	1/1	0.99	0.09	1,1,1,1	0
57	MG	DE	301	1/1	0.99	0.20	1,1,1,1	1
57	MG	CA	1800	1/1	0.99	0.09	9,9,9,9	0
57	MG	DA	3125	1/1	0.99	0.39	4,4,4,4	1
57	MG	BA	3096	1/1	0.99	0.32	1,1,1,1	1
57	MG	BA	3319	1/1	0.99	0.11	1,1,1,1	0
57	MG	DA	3083	1/1	0.99	0.23	22,22,22,22	0
57	MG	BA	3016	1/1	0.99	0.14	31,31,31,31	1
57	MG	AA	1730	1/1	0.99	0.13	1,1,1,1	1
57	MG	BB	205	1/1	0.99	0.10	6,6,6,6	1
57	MG	DA	3267	1/1	0.99	0.24	28,28,28,28	1
57	MG	CA	1618	1/1	0.99	0.07	22,22,22,22	1
57	MG	DA	3342	1/1	0.99	0.22	1,1,1,1	1
57	MG	DA	3073	1/1	0.99	0.21	1,1,1,1	1
57	MG	CA	1795	1/1	0.99	0.14	34,34,34,34	0
57	MG	BA	3031	1/1	0.99	0.30	1,1,1,1	0
57	MG	AA	1649	1/1	0.99	0.09	56,56,56,56	0
57	MG	BA	3170	1/1	0.99	0.08	1,1,1,1	1
57	MG	BA	3007	1/1	0.99	0.09	37,37,37,37	0
57	MG	BA	3069	1/1	0.99	0.27	1,1,1,1	0
57	MG	BA	3381	1/1	0.99	0.11	23,23,23,23	0
57	MG	BA	3212	1/1	0.99	0.06	2,2,2,2	0
57	MG	DA	3239	1/1	0.99	0.14	4,4,4,4	0
57	MG	DA	3072	1/1	0.99	0.09	1,1,1,1	1
57	MG	DA	3078	1/1	0.99	0.16	58,58,58,58	1
57	MG	DA	3182	1/1	0.99	0.19	1,1,1,1	1
57	MG	D1	103	1/1	0.99	0.08	33,33,33,33	0
57	MG	BA	3162	1/1	0.99	0.22	54,54,54,54	1
57	MG	DA	3122	1/1	0.99	0.19	8,8,8,8	0
57	MG	BA	3068	1/1	0.99	0.20	1,1,1,1	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3298	1/1	0.99	0.17	1,1,1,1	1
57	MG	BA	3247	1/1	0.99	0.27	26,26,26,26	1
57	MG	DA	3064	1/1	0.99	0.09	18,18,18,18	0
57	MG	BA	3057	1/1	0.99	0.16	7,7,7,7	0
57	MG	AA	1711	1/1	0.99	0.14	21,21,21,21	1
57	MG	CA	1605	1/1	0.99	0.08	18,18,18,18	0
57	MG	BA	3124	1/1	0.99	0.09	7,7,7,7	0
57	MG	BB	216	1/1	0.99	0.35	1,1,1,1	1
57	MG	CA	1609	1/1	0.99	0.21	62,62,62,62	0
57	MG	DA	3020	1/1	0.99	0.48	1,1,1,1	0
57	MG	DA	3234	1/1	0.99	0.11	88,88,88,88	1
57	MG	BA	3078	1/1	0.99	0.17	1,1,1,1	0
57	MG	BA	3262	1/1	0.99	0.18	32,32,32,32	1
57	MG	DA	3044	1/1	0.99	0.27	60,60,60,60	0
57	MG	BA	3013	1/1	0.99	0.28	1,1,1,1	0
57	MG	BA	3017	1/1	0.99	0.12	3,3,3,3	1
57	MG	CA	1601	1/1	0.99	0.12	21,21,21,21	1
57	MG	BA	3161	1/1	0.99	0.19	7,7,7,7	1
57	MG	BA	3244	1/1	0.99	0.14	1,1,1,1	0
57	MG	BA	3220	1/1	0.99	0.16	16,16,16,16	1
57	MG	AA	1783	1/1	0.99	0.09	39,39,39,39	1
57	MG	DA	3120	1/1	0.99	0.19	1,1,1,1	1
57	MG	BA	3283	1/1	0.99	0.16	2,2,2,2	0
57	MG	BA	3134	1/1	0.99	0.07	1,1,1,1	0
57	MG	BA	3050	1/1	0.99	0.14	1,1,1,1	1
57	MG	DA	3172	1/1	0.99	0.15	17,17,17,17	1
57	MG	DA	3030	1/1	0.99	0.19	2,2,2,2	0
57	MG	DA	3347	1/1	0.99	0.32	35,35,35,35	1
57	MG	DA	3079	1/1	0.99	0.09	5,5,5,5	0
57	MG	DA	3185	1/1	0.99	0.16	19,19,19,19	0
57	MG	DA	3080	1/1	0.99	0.09	1,1,1,1	1
57	MG	AV	104	1/1	0.99	0.08	58,58,58,58	1
57	MG	AA	1620	1/1	0.99	0.17	3,3,3,3	0
57	MG	AA	1689	1/1	0.99	0.14	6,6,6,6	0
57	MG	BA	3074	1/1	0.99	0.08	1,1,1,1	1
57	MG	DA	3123	1/1	0.99	0.22	1,1,1,1	0
57	MG	DA	3257	1/1	0.99	0.15	54,54,54,54	0
57	MG	DD	301	1/1	0.99	0.19	1,1,1,1	0
57	MG	DA	3235	1/1	0.99	0.12	7,7,7,7	1
57	MG	DA	3019	1/1	0.99	0.10	23,23,23,23	0
57	MG	DA	3011	1/1	0.99	0.11	42,42,42,42	0
57	MG	DA	3038	1/1	0.99	0.20	3,3,3,3	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
57	MG	BA	3052	1/1	0.99	0.10	51,51,51,51	0
57	MG	AA	1622	1/1	0.99	0.08	1,1,1,1	1
57	MG	AA	1604	1/1	0.99	0.08	3,3,3,3	0
57	MG	DA	3121	1/1	0.99	0.37	20,20,20,20	0
57	MG	BA	3095	1/1	0.99	0.12	1,1,1,1	0
57	MG	CA	1696	1/1	0.99	0.66	52,52,52,52	0
57	MG	DA	3259	1/1	0.99	0.18	22,22,22,22	0
57	MG	DA	3378	1/1	0.99	0.21	18,18,18,18	1
57	MG	DA	3417	1/1	1.00	0.10	71,71,71,71	0
58	ZN	AD	303	1/1	1.00	0.20	26,26,26,26	0
57	MG	DA	3208	1/1	1.00	0.11	14,14,14,14	0
58	ZN	CD	301	1/1	1.00	0.22	35,35,35,35	0
57	MG	BA	3436	1/1	1.00	0.18	6,6,6,6	1

6.5 Other polymers [i](#)

There are no such residues in this entry.