



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 03:47 pm BST

PDB ID : 4V56
Title : Crystal structure of the bacterial ribosome from Escherichia coli in complex with spectinomycin.
Authors : Borovinskaya, M.A.; Shoji, S.; Holton, J.M.; Fredrick, K.; Cate, J.H.D.
Deposited on : 2007-07-21
Resolution : 3.93 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
buster-report : 1.1.7 (2018)
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

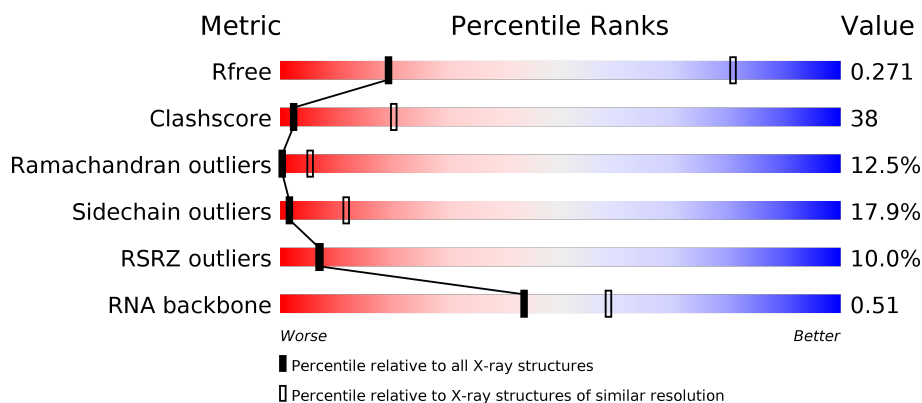
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.93 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	1036 (4.20-3.68)
Clashscore	141614	1009 (4.18-3.70)
Ramachandran outliers	138981	1057 (4.20-3.68)
Sidechain outliers	138945	1049 (4.20-3.68)
RSRZ outliers	127900	1007 (4.24-3.64)
RNA backbone	3102	1041 (4.84-3.00)

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	1542	<div> <div>19%</div> <div>63%</div> <div>16%</div> <div>..</div> </div>
1	CA	1542	<div> <div>20%</div> <div>63%</div> <div>16%</div> <div>.</div> </div>
2	AC	232	<div> <div>6%</div> <div>19%</div> <div>46%</div> <div>22%</div> <div>11%</div> <div>.</div> </div>
2	CC	232	<div> <div>12%</div> <div>22%</div> <div>48%</div> <div>18%</div> <div>11%</div> </div>

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Mol	Chain	Length	Quality of chain
3	AD	205	
3	CD	205	
4	AE	166	
4	CE	166	
5	AF	135	
5	CF	135	
6	AG	178	
6	CG	178	
7	AH	129	
7	CH	129	
8	AI	129	
8	CI	129	
9	AJ	103	
9	CJ	103	
10	AK	128	
10	CK	128	
11	AL	123	
11	CL	123	
12	AM	117	
12	CM	117	
13	AP	82	
13	CP	82	
14	AQ	83	
14	CQ	83	
15	AR	74	

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Mol	Chain	Length	Quality of chain
15	CR	74	
16	AS	91	
16	CS	91	
17	AT	86	
17	CT	86	
18	AB	240	
18	CB	240	
19	AU	70	
19	CU	70	
20	AO	89	
20	CO	89	
21	AN	100	
21	CN	100	
22	BA	120	
22	DA	120	
23	BB	2904	
23	DB	2904	
24	BI	141	
24	DI	141	
25	BC	272	
25	DC	272	
26	BD	209	
26	DD	209	
27	BK	123	
27	DK	123	

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Mol	Chain	Length	Quality of chain
28	BP	114	
28	DP	114	
29	BE	201	
29	DE	201	
30	BY	58	
30	DY	58	
31	B0	56	
31	D0	56	
32	B4	38	
32	D4	38	
33	B1	54	
33	D1	54	
34	B3	64	
34	D3	64	
35	BV	94	
35	DV	94	
36	B2	46	
36	D2	46	
37	BL	144	
37	DL	144	
38	BM	136	
38	DM	136	
39	BX	63	
39	DX	63	
40	BH	149	

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Mol	Chain	Length	Quality of chain
40	DH	149	
41	BJ	142	
41	DJ	142	
42	BN	127	
42	DN	127	
43	BO	117	
43	DO	117	
44	BQ	117	
44	DQ	117	
45	BS	110	
45	DS	110	
46	BU	103	
46	DU	103	
47	BF	178	
47	DF	178	
48	BG	176	
48	DG	176	
49	BR	103	
49	DR	103	
50	BT	100	
50	DT	100	
51	BZ	78	
51	DZ	78	
52	BW	84	
52	DW	84	

2 Entry composition

There are 56 unique types of molecules in this entry. The entry contains 284033 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 rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	1530	Total	C	N	O	P	0	0	0
			32831	14642	6024	10635	1530			
1	CA	1530	Total	C	N	O	P	0	0	0
			32831	14642	6024	10635	1530			

- Molecule 2 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AC	206	Total	C	N	O	S	0	0	0
			1624	1028	305	288	3			
2	CC	206	Total	C	N	O	S	0	0	0
			1624	1028	305	288	3			

- Molecule 3 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AD	205	Total	C	N	O	S	0	0	0
			1643	1026	315	298	4			
3	CD	205	Total	C	N	O	S	0	0	0
			1643	1026	315	298	4			

- Molecule 4 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AE	150	Total	C	N	O	S	0	0	0
			1105	687	211	201	6			
4	CE	150	Total	C	N	O	S	0	0	0
			1105	687	211	201	6			

- Molecule 5 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AF	100	Total	C	N	O	S	0	0	0
			817	515	148	148	6			
5	CF	100	Total	C	N	O	S	0	0	0
			817	515	148	148	6			

- Molecule 6 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AG	150	Total	C	N	O	S	0	0	0
			1174	730	226	214	4			
6	CG	152	Total	C	N	O	S	0	0	0
			1196	745	230	217	4			

- Molecule 7 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AH	129	Total	C	N	O	S	0	0	0
			979	616	173	184	6			
7	CH	129	Total	C	N	O	S	0	0	0
			979	616	173	184	6			

- Molecule 8 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AI	127	Total	C	N	O	S	0	0	0
			1022	634	206	179	3			
8	CI	127	Total	C	N	O	S	0	0	0
			1022	634	206	179	3			

- Molecule 9 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	AJ	98	Total	C	N	O	S	0	0	0
			786	493	150	142	1			
9	CJ	98	Total	C	N	O	S	0	0	0
			786	493	150	142	1			

- Molecule 10 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AK	117	Total	C	N	O	S	0	0	0
			877	540	174	160	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	CK	117	Total	C	N	O	S	0	0	0
			877	540	174	160	3			

- Molecule 11 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AL	123	Total	C	N	O	S	0	0	0
			955	590	196	165	4			
11	CL	123	Total	C	N	O	S	0	0	0
			955	590	196	165	4			

- Molecule 12 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AM	114	Total	C	N	O	S	0	0	0
			883	546	178	156	3			
12	CM	113	Total	C	N	O	S	0	0	0
			876	541	177	155	3			

- Molecule 13 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AP	82	Total	C	N	O	S	0	0	0
			649	406	128	114	1			
13	CP	80	Total	C	N	O	S	0	0	0
			638	400	126	111	1			

- Molecule 14 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AQ	80	Total	C	N	O	S	0	0	0
			648	411	121	113	3			
14	CQ	81	Total	C	N	O	S	0	0	0
			657	417	122	115	3			

- Molecule 15 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	AR	55	Total	C	N	O	0	0	0
			455	288	86	81			
15	CR	55	Total	C	N	O	0	0	0
			455	288	86	81			

- Molecule 16 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AS	79	Total	C	N	O	S	0	0	0
			637	408	120	107	2			
16	CS	80	Total	C	N	O	S	0	0	0
			644	413	121	108	2			

- Molecule 17 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AT	85	Total	C	N	O	S	0	0	0
			665	411	137	114	3			
17	CT	85	Total	C	N	O	S	0	0	0
			665	411	137	114	3			

- Molecule 18 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	AB	218	Total	C	N	O	S	0	0	0
			1704	1081	305	311	7			
18	CB	218	Total	C	N	O	S	0	0	0
			1704	1081	305	311	7			

- Molecule 19 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AU	51	Total	C	N	O	S	0	0	0
			425	265	86	73	1			
19	CU	51	Total	C	N	O	S	0	0	0
			425	265	86	73	1			

- Molecule 20 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AO	88	Total	C	N	O	S	0	0	0
			714	439	144	130	1			
20	CO	88	Total	C	N	O	S	0	0	0
			714	439	144	130	1			

- Molecule 21 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	AN	96	Total	C	N	O	S	0	0	0
			774	483	160	128	3			
21	CN	96	Total	C	N	O	S	0	0	0
			774	483	160	128	3			

- Molecule 22 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	BA	117	Total	C	N	O	P	0	0	0
			2507	1116	459	815	117			
22	DA	117	Total	C	N	O	P	0	0	0
			2507	1116	459	815	117			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BA	120	U	-	INSERTION	GB 85674274
DA	120	U	-	INSERTION	GB 85674274

- Molecule 23 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	BB	2841	Total	C	N	O	P	0	0	0
			60995	27210	11229	19715	2841			
23	DB	2841	Total	C	N	O	P	0	0	0
			60995	27210	11229	19715	2841			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BB	2903	U	-	INSERTION	GB 85674274
BB	2904	U	-	INSERTION	GB 85674274
DB	2903	U	-	INSERTION	GB 85674274
DB	2904	U	-	INSERTION	GB 85674274

- Molecule 24 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	BI	141	Total	C	N	O	S	0	0	0
			1032	651	179	196	6			
24	DI	141	Total	C	N	O	S	0	0	0
			1032	651	179	196	6			

- Molecule 25 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	BC	271	Total	C	N	O	S	0	0	0
			2082	1288	423	364	7			
25	DC	271	Total	C	N	O	S	0	0	0
			2082	1288	423	364	7			

- Molecule 26 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	BD	209	Total	C	N	O	S	0	0	0
			1565	979	288	294	4			
26	DD	209	Total	C	N	O	S	0	0	0
			1565	979	288	294	4			

- Molecule 27 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	BK	121	Total	C	N	O	S	0	0	0
			930	582	179	164	5			
27	DK	121	Total	C	N	O	S	0	0	0
			930	582	179	164	5			

- Molecule 28 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	BP	114	Total	C	N	O	S	0	0	0
			917	574	179	163	1			
28	DP	114	Total	C	N	O	S	0	0	0
			917	574	179	163	1			

- Molecule 29 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	BE	201	Total	C	N	O	S	0	0	0
			1552	974	283	290	5			
29	DE	201	Total	C	N	O	S	0	0	0
			1552	974	283	290	5			

- Molecule 30 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	BY	58	Total	C	N	O	S	0	0	0
			449	281	87	79	2			
30	DY	58	Total	C	N	O	S	0	0	0
			449	281	87	79	2			

- Molecule 31 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	B0	56	Total	C	N	O	S	0	0	0
			444	269	94	80	1			
31	D0	56	Total	C	N	O	S	0	0	0
			444	269	94	80	1			

- Molecule 32 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	B4	38	Total	C	N	O	S	0	0	0
			302	185	65	48	4			
32	D4	38	Total	C	N	O	S	0	0	0
			302	185	65	48	4			

- Molecule 33 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
33	B1	50	Total	C	N	O	0	0	0
			409	263	75	71			
33	D1	50	Total	C	N	O	0	0	0
			409	263	75	71			

- Molecule 34 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	B3	64	Total	C	N	O	S	0	0	0
			504	323	105	74	2			
34	D3	64	Total	C	N	O	S	0	0	0
			504	323	105	74	2			

- Molecule 35 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BV	94	Total	C	N	O	S	0	0	0
			753	479	137	134	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	DV	94	Total	C	N	O	S	0	0	0
			753	479	137	134	3			

- Molecule 36 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	B2	46	Total	C	N	O	S	0	0	0
			377	228	90	57	2			
36	D2	46	Total	C	N	O	S	0	0	0
			377	228	90	57	2			

- Molecule 37 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	BL	143	Total	C	N	O	S	0	0	0
			1045	649	206	189	1			
37	DL	143	Total	C	N	O	S	0	0	0
			1045	649	206	189	1			

- Molecule 38 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	BM	136	Total	C	N	O	S	0	0	0
			1074	686	205	177	6			
38	DM	136	Total	C	N	O	S	0	0	0
			1074	686	205	177	6			

- Molecule 39 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	BX	63	Total	C	N	O	S	0	0	0
			509	313	99	95	2			
39	DX	63	Total	C	N	O	S	0	0	0
			509	313	99	95	2			

- Molecule 40 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	BH	149	Total	C	N	O	S	0	0	0
			1111	699	197	214	1			
40	DH	149	Total	C	N	O	S	0	0	0
			1111	699	197	214	1			

- Molecule 41 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BJ	142	Total	C	N	O	S	0	0	0
			1129	714	212	199	4			
41	DJ	142	Total	C	N	O	S	0	0	0
			1129	714	212	199	4			

- Molecule 42 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BN	120	Total	C	N	O	S	0	0	0
			960	593	196	166	5			
42	DN	120	Total	C	N	O	S	0	0	0
			960	593	196	166	5			

- Molecule 43 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
43	BO	116	Total	C	N	O	0	0	0
			892	552	178	162			
43	DO	116	Total	C	N	O	0	0	0
			892	552	178	162			

- Molecule 44 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
44	BQ	117	Total	C	N	O	0	0	0
			947	604	192	151			
44	DQ	117	Total	C	N	O	0	0	0
			947	604	192	151			

- Molecule 45 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	BS	110	Total	C	N	O	S	0	0	0
			857	532	166	156	3			
45	DS	110	Total	C	N	O	S	0	0	0
			857	532	166	156	3			

- Molecule 46 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
46	BU	102	Total	C	N	O	0	0	0
			779	492	146	141			
46	DU	102	Total	C	N	O	0	0	0
			779	492	146	141			

- Molecule 47 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	BF	178	Total	C	N	O	S	0	0	0
			1420	905	251	258	6			
47	DF	178	Total	C	N	O	S	0	0	0
			1420	905	251	258	6			

- Molecule 48 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	BG	176	Total	C	N	O	S	0	0	0
			1323	832	243	246	2			
48	DG	176	Total	C	N	O	S	0	0	0
			1323	832	243	246	2			

- Molecule 49 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	BR	103	Total	C	N	O	S	0	0	0
			816	516	153	145	2			
49	DR	103	Total	C	N	O	S	0	0	0
			816	516	153	145	2			

- Molecule 50 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	BT	93	Total	C	N	O	S	0	0	0
			738	466	139	131	2			
50	DT	93	Total	C	N	O	S	0	0	0
			738	466	139	131	2			

- Molecule 51 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	BZ	77	Total	C	N	O	S	0	0	0
			625	388	129	106	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	DZ	77	Total	C	N	O	S	0	0	0
			625	388	129	106	2			

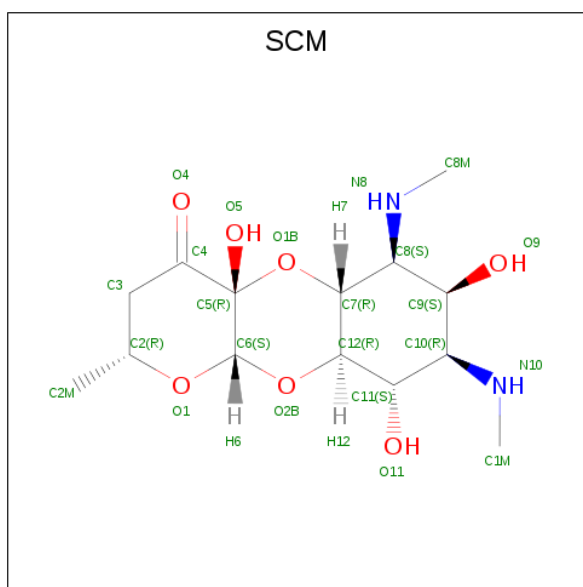
- Molecule 52 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	BW	79	Total	C	N	O	S	0	0	0
			596	367	120	108	1			
52	DW	79	Total	C	N	O	S	0	0	0
			596	367	120	108	1			

- Molecule 53 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
53	BB	110	Total	Mg	0	0
			110	110		
53	CA	58	Total	Mg	0	0
			58	58		
53	AA	60	Total	Mg	0	0
			60	60		
53	CE	1	Total	Mg	0	0
			1	1		
53	DN	1	Total	Mg	0	0
			1	1		
53	DB	110	Total	Mg	0	0
			110	110		

- Molecule 54 is SPECTINOMYCIN (three-letter code: SCM) (formula: C₁₄H₂₄N₂O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
54	AA	1	Total	C	N	O	0	0
			23	14	2	7		
54	CA	1	Total	C	N	O	0	0
			23	14	2	7		

- Molecule 55 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
55	B4	1	Total	Zn	0	0
			1	1		
55	D4	1	Total	Zn	0	0
			1	1		

- Molecule 56 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	AA	288	Total	O	0	0
			288	288		
56	AE	3	Total	O	0	0
			3	3		
56	AK	1	Total	O	0	0
			1	1		
56	AL	4	Total	O	0	0
			4	4		
56	AP	1	Total	O	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	AT	1	Total 1	O 1	0	0
56	AN	2	Total 2	O 2	0	0
56	BB	494	Total 494	O 494	0	0
56	BC	4	Total 4	O 4	0	0
56	BE	3	Total 3	O 3	0	0
56	BL	4	Total 4	O 4	0	0
56	BH	1	Total 1	O 1	0	0
56	BT	1	Total 1	O 1	0	0
56	CA	275	Total 275	O 275	0	0
56	CE	4	Total 4	O 4	0	0
56	CK	1	Total 1	O 1	0	0
56	CL	5	Total 5	O 5	0	0
56	CP	1	Total 1	O 1	0	0
56	CT	2	Total 2	O 2	0	0
56	CN	5	Total 5	O 5	0	0
56	DB	500	Total 500	O 500	0	0
56	DC	3	Total 3	O 3	0	0
56	DD	1	Total 1	O 1	0	0
56	DP	1	Total 1	O 1	0	0
56	DE	1	Total 1	O 1	0	0
56	DL	3	Total 3	O 3	0	0

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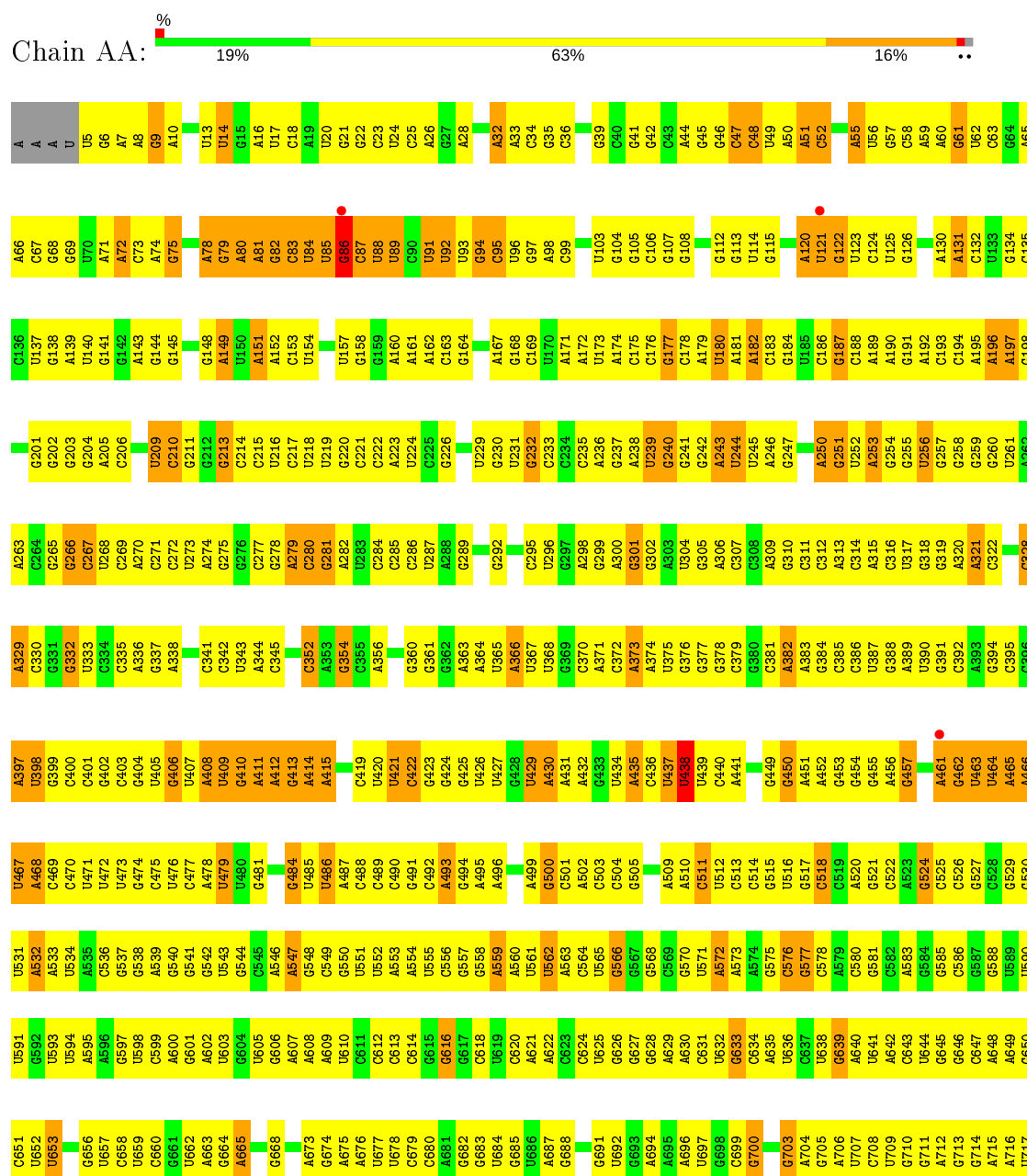
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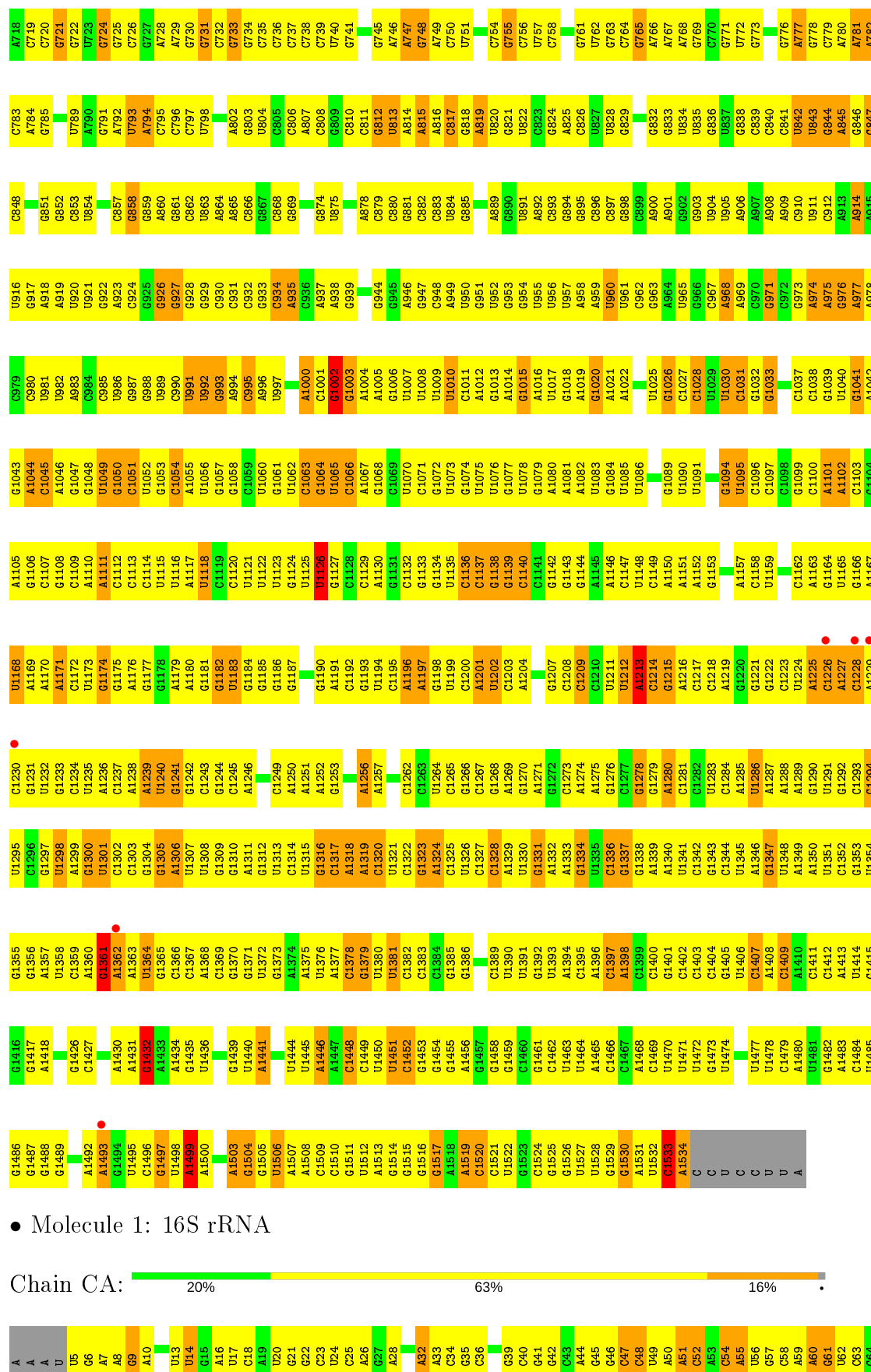
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	DJ	1	Total	O	0	0
			1	1		
56	DN	2	Total	O	0	0
			2	2		
56	DR	1	Total	O	0	0
			1	1		

3 Residue-property plots

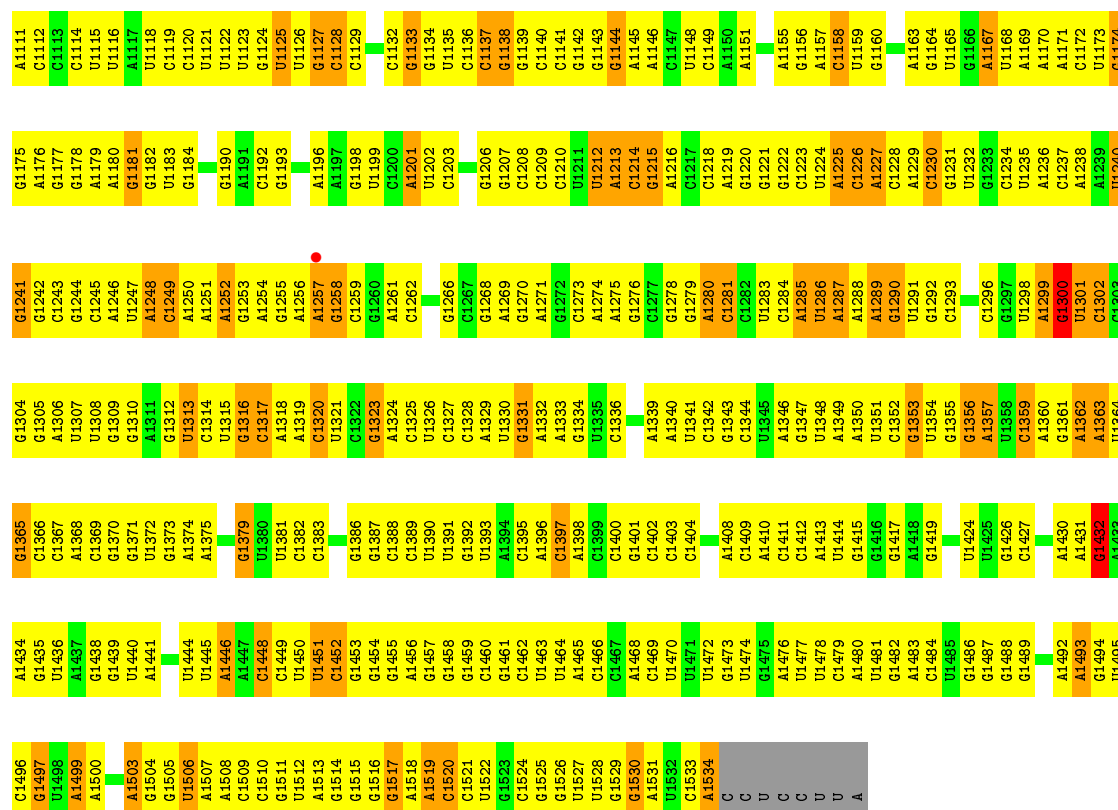
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 rRNA

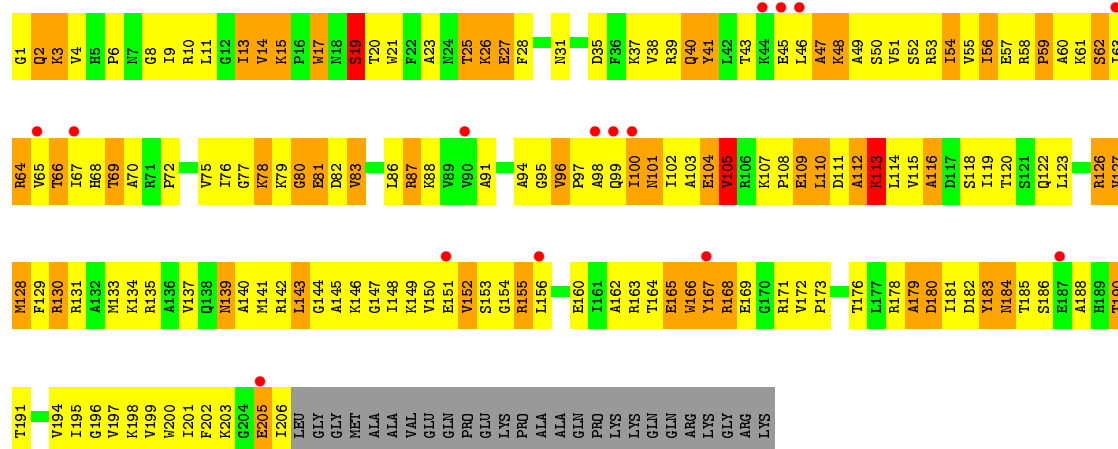
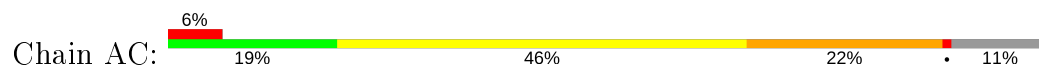




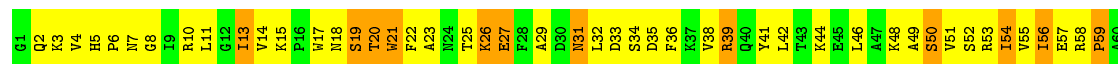
G1048	G1049	G985	G922	C857	A792	G724	C660	G597	G537	G474	G402	C335	G266	G203	U133	A65
U986	U987	G858	A923	G856	U793	G725	G661	U598	G538	C475	C403	A336	C267	G204	G134	A66
G988	G989	A860	C924	A859	A794	G728	U662	C599	A539	U476	U405	A337	U268	C206	C136	G67
U989	U990	G861	G925	A860	C795	A728	A663	A600	G540	U477	U406	A338	C269	G206	G137	G68
G990	G991	C862	G926	C861	C796	A729	G664	G601	G541	A478	G407	C341	A270	U209	G138	U70
C1054	C1055	U863	G928	U863	U798	G731	G666	U603	U543	U480	A408	C342	C271	C210	A143	A71
A1056	A1057	A864	G929	A864	C732	G733	G667	G604	C544	G481	U409	U343	U272	C212	A144	A72
U1056	U1058	A865	C930	C866	G733	G734	G668	U605	C545	A482	G410	U344	U273	G211	G144	C73
G1057	G1058	C866	G931	G867	C734	G735	A483	G606	C483	C483	A411	C345	A274	G212	G145	A74
A994	A995	G867	C932	G867	U804	C735	U672	A607	A547	A484	A412	C346	G275	C214	G145	G75
A996	A997	C868	G933	C868	G736	C736	A673	A608	G548	U485	G413	G347	A279	C215	G148	G76
U997	U998	G869	C934	G869	C737	C737	G674	A609	C549	U486	A414	C352	G280	C216	A149	A77
G1061	G1062	A807	A935	A807	C738	C738	A675	G609	G550	A487	A415	C353	G281	C217	U150	A78
A1000	A1001	C808	G936	C808	C739	C739	A676	C612	U551	C488	U420	A354	A282	U218	U151	G79
C1001	C1002	U875	A937	U875	U740	U740	A677	C613	U552	C489	U421	U345	U283	U219	A152	A80
G1003	G1004	G881	A938	C876	C810	G741	U678	C614	A553	C490	U422	C355	C284	G220	C153	A81
U1005	U1006	C811	G939	G877	C810	C679	U679	G615	A554	A491	G423	A356	C285	G221	U154	G82
A1004	A1005	G812	C940	A878	C680	C745	C680	G616	U555	C492	G424	C361	C286	C222	C83	C83
C1066	C1067	U813	G941	C879	A681	A746	A681	G617	C556	A493	G425	G362	U287	C223	U157	U84
A1008	A1009	A814	A942	C880	G682	A747	A682	C618	G557	G494	G426	A363	U288	U224	G158	U85
G1068	G1069	A815	G943	G881	U683	G748	U683	U619	G558	A495	U427	A364	G289	C225	G159	G86
C1007	C1008	A816	C882	C882	U684	A749	A559	C620	G559	A496	U428	U365	G300	G226	A160	C87
U1070	U1071	C817	G944	C883	G685	C750	G685	A621	A560	U429	G429	U366	G301	C227	A161	U88
G1072	G1073	G818	C948	U884	U686	G753	U686	A622	U561	A499	A430	U367	G302	U229	A162	U89
U1074	U1075	A819	G949	G885	A687	A753	A687	C623	U562	G500	A431	U368	G303	G230	C163	C90
G1076	G1077	U820	U950	C889	G688	C754	G688	C624	A563	C501	A432	U369	G304	U231	G164	U91
U1078	U1079	G821	G951	A889	G689	G755	G689	U625	A564	A502	A433	G370	U304	G232	U92	U92
A1080	A1081	U822	G952	C890	U691	C756	U691	G633	U571	U339	U438	G371	U305	C233	A167	U93
C1081	C1082	C823	U891	G890	U692	U757	U692	C634	G504	A434	U434	A371	G299	C234	G178	G102
U1083	U1084	G824	U892	A892	G693	C758	G693	A629	G505	A435	A435	C372	G300	C235	C169	G94
G1085	G1086	A825	C893	C893	A694	C759	A694	A630	G568	A509	G436	A373	G301	A236	G179	C95
U1088	U1089	G826	G894	A895	A695	A761	A695	C631	G569	A510	U437	A374	G302	G237	A174	A98
A1022	A1023	U827	C895	C895	U696	G762	U696	U632	G570	U375	U438	G376	G303	A238	C176	C99
U1024	U1025	G829	C897	C896	U697	G763	U697	G633	G571	C511	U439	G377	G304	U239	G177	G100
G1026	G1027	U835	G902	C897	G698	C764	G698	C634	A572	U512	G441	G378	G305	G240	G178	A101
C1028	C1029	G836	U904	G903	C699	G765	C699	A635	A573	C513	G449	G379	G306	G241	G103	G102
U1029	U1030	U837	U905	U904	G700	A766	G700	U636	A574	C514	G450	C379	G307	G242	U103	U103
G1031	G1032	G838	U906	U906	U707	A767	U707	G637	G575	G515	G451	G380	G308	A243	U180	G104
U1033	U1034	C839	A906	A906	C708	A767	A702	U638	G576	U516	A451	C381	A309	U244	A181	G105
C1035	C1036	C840	A907	A907	U709	A768	G703	U639	G577	G517	A452	A382	G310	U245	A182	C106
A1035	A1036	C841	A908	A909	G710	G776	G703	A640	C578	G518	G453	A383	G311	A246	C183	G107
G1037	G1038	U842	G909	G909	G711	G777	A704	G646	G584	C519	G454	G384	G312	G247	U184	G108
C1039	C1040	U843	C910	C910	G712	G778	G705	U641	G585	A520	G455	G385	G313	G248	U185	G109
U1041	U1042	G844	U911	U911	A712	C779	G706	A642	G586	G521	A456	C386	G314	G249	A186	G112
A1042	A1043	A845	G912	C912	G713	A780	A706	U643	G587	G522	G457	C387	G315	A250	C186	G113
C1043	C1044	G846	A913	A913	G714	G781	C708	U644	G588	A523	A461	U387	A316	G251	G187	U114
G1045	G1046	U847	A914	A914	G715	A782	U709	G645	A583	G524	G462	U388	G317	U252	C188	G115
A1046	A1047	C848	A915	A915	A716	C783	G710	G650	G584	G525	A463	U389	U317	A253	A190	A120
G1047	G1048	A717	G916	G916	G717	A784	G711	U652	G585	C526	U464	U390	G318	G254	G191	U121
C1049	C1050	U849	U917	U917	A718	G785	G712	A648	G586	G527	U465	U391	G319	U255	A192	U122
U1051	U1052	G851	G917	G917	G719	G786	G713	A649	G587	G528	A466	G392	A321	U256	A193	U123
A1052	A1053	C852	A918	A918	G720	G787	G714	G650	G588	C529	A467	A393	C322	G257	C193	U124
G1054	G1055	U854	A919	A919	G721	U789	A715	C651	U589	G529	U467	U394	G328	G258	C194	U125
C1056	C1057	G855	U920	U920	G722	A790	A716	U652	U590	G530	A468	A395	A329	G259	A195	C124
U1058	U1059	A915	G918	G918	G723	G791	G717	U653	U591	U531	C469	C396	A330	U261	A196	U126
G1060	G1061	G856	A919	A919	G724	G792	G718	G656	G592	A532	C470	U397	G331	U262	A197	G126
A1062	A1063	U857	U921	U921	G725	G793	G719	U657	U593	A533	C471	U398	G332	G263	G198	A130
C1064	C1065	G858	G920	G920	G726	U794	G720	U658	U594	A534	U471	U399	G333	C264	G201	A131
U1066	U1067	A915	A921	A921	G727	G795	G721	G659	U595	A535	U472	C400	U333	G265	G202	C132
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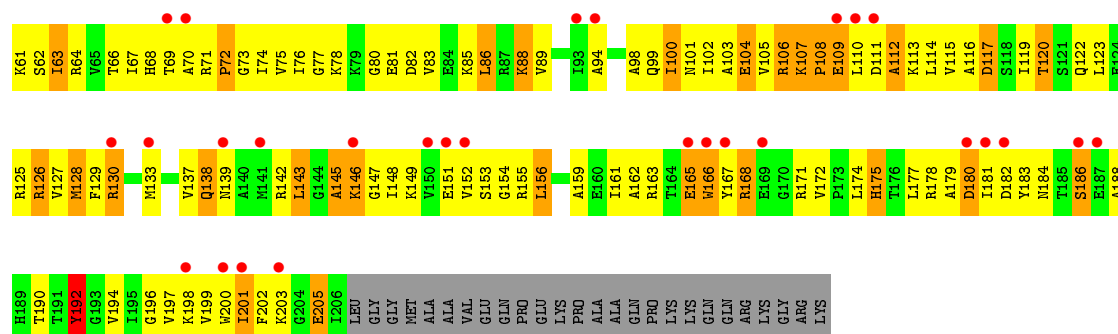


• Molecule 2: 30S ribosomal protein S3

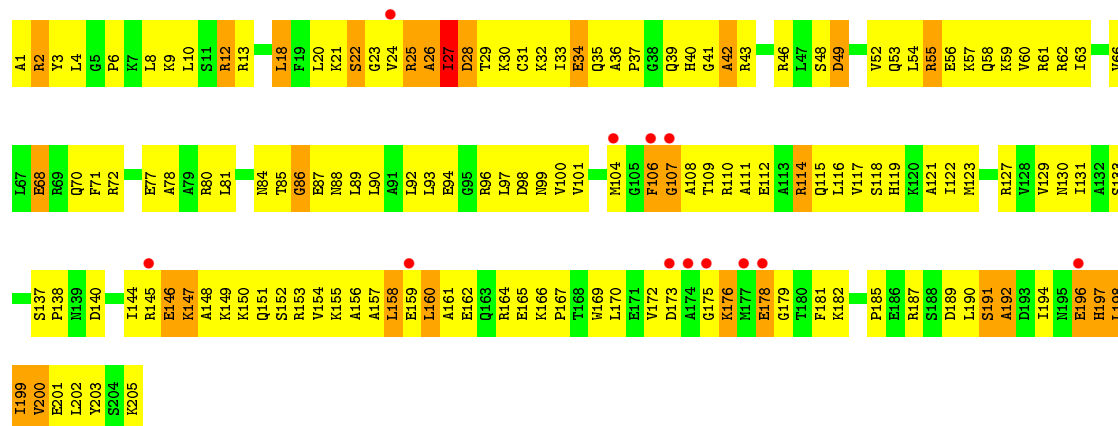


• Molecule 2: 30S ribosomal protein S3

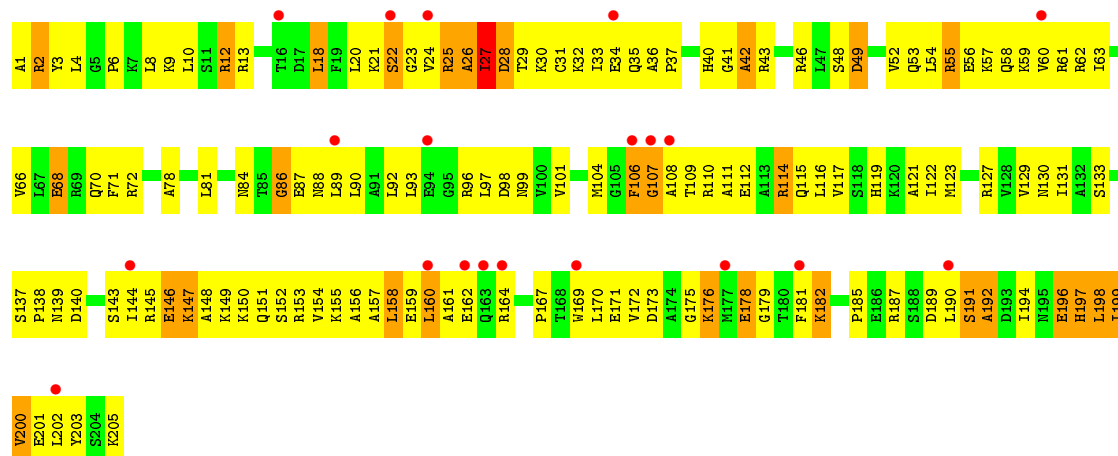




• Molecule 3: 30S ribosomal protein S4

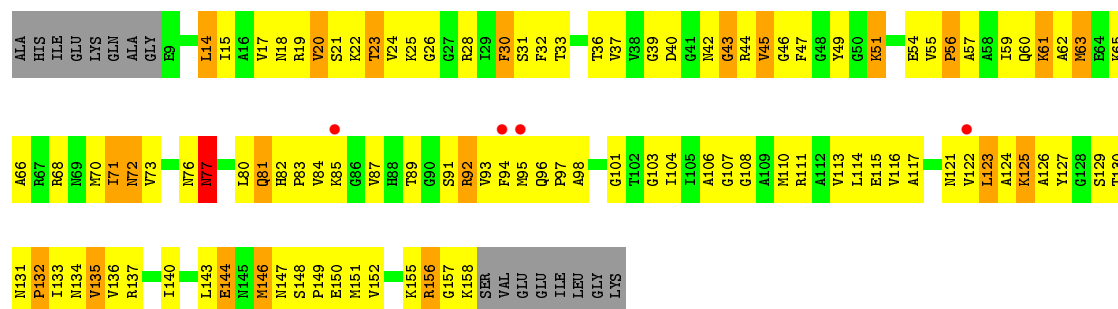


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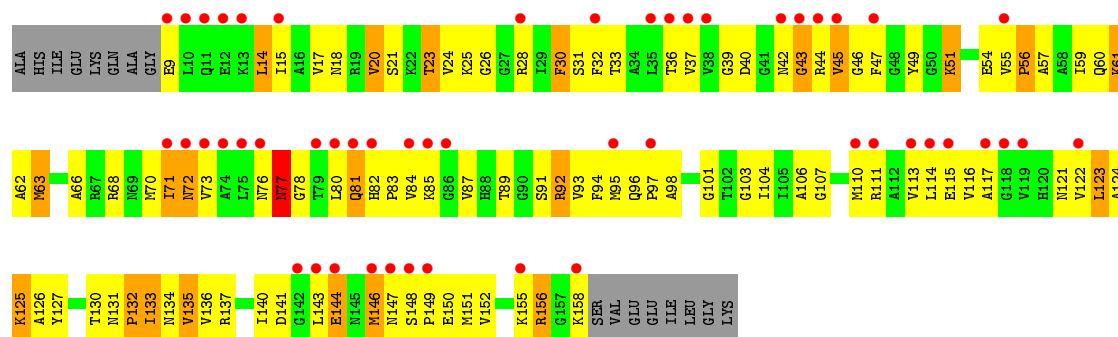


• Molecule 4: 30S ribosomal protein S5

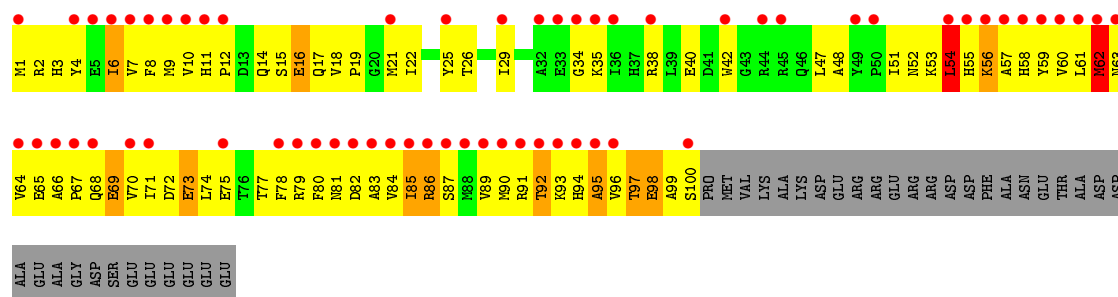
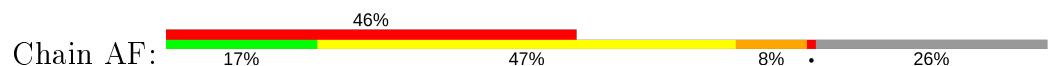




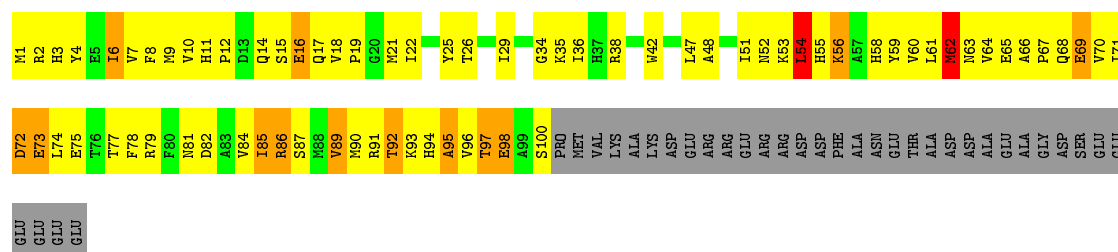
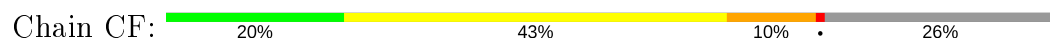
• Molecule 4: 30S ribosomal protein S5



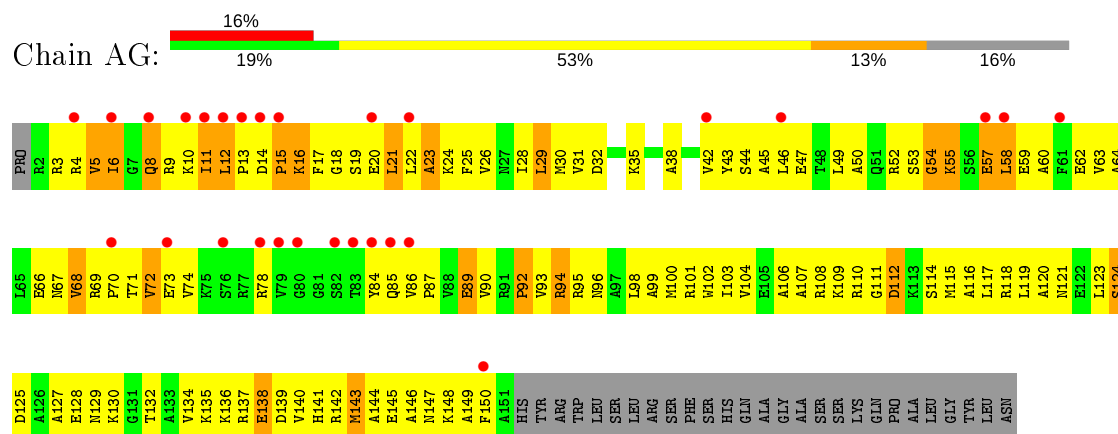
• Molecule 5: 30S ribosomal protein S6



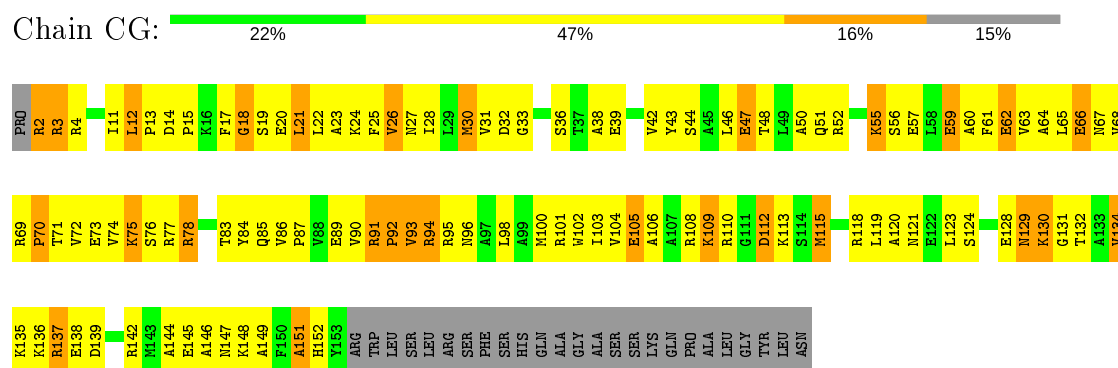
• Molecule 5: 30S ribosomal protein S6



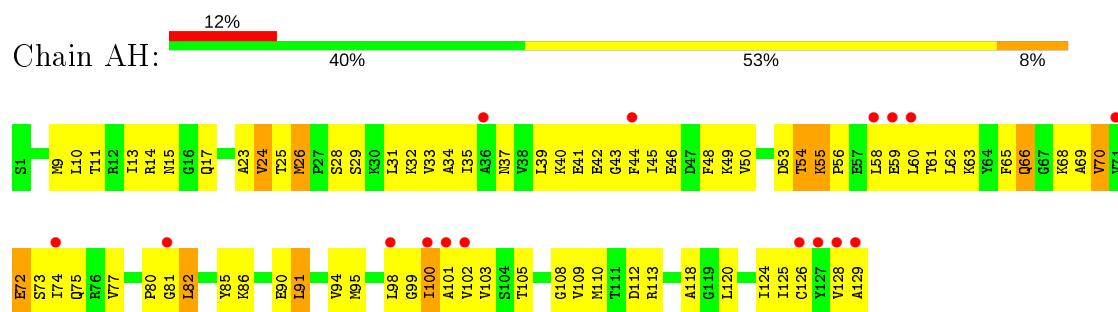
- Molecule 6: 30S ribosomal protein S7



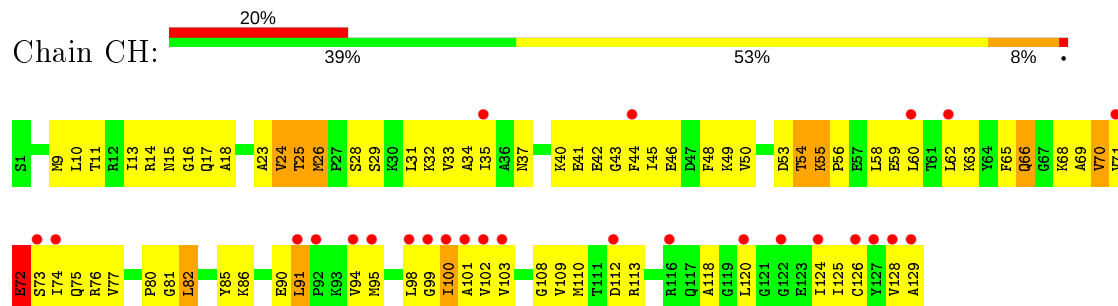
- Molecule 6: 30S ribosomal protein S7



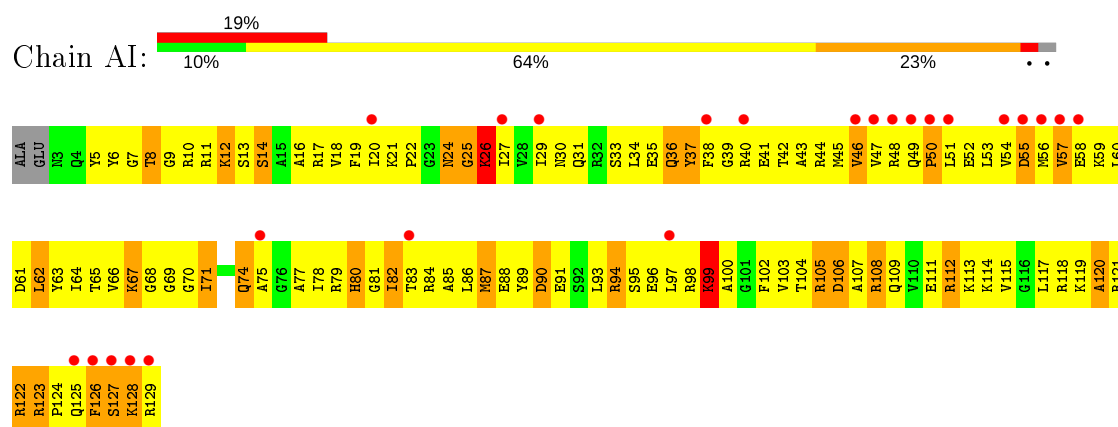
- Molecule 7: 30S ribosomal protein S8



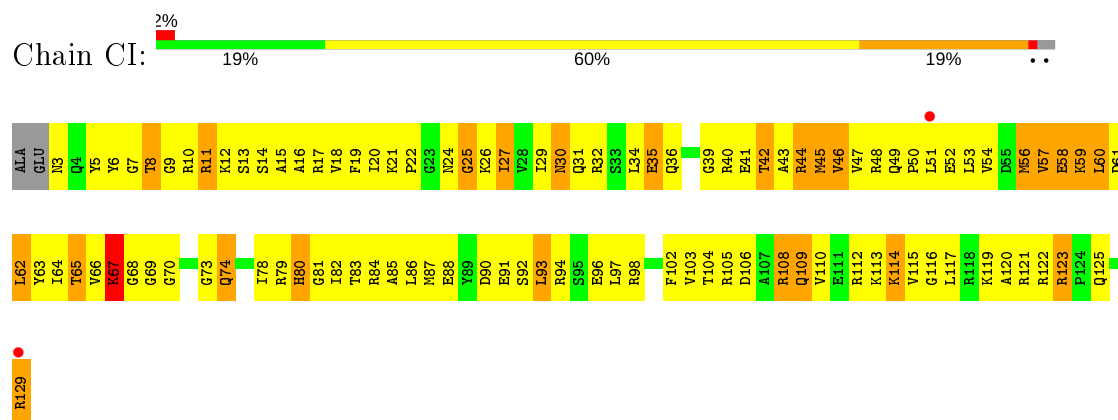
- Molecule 7: 30S ribosomal protein S8



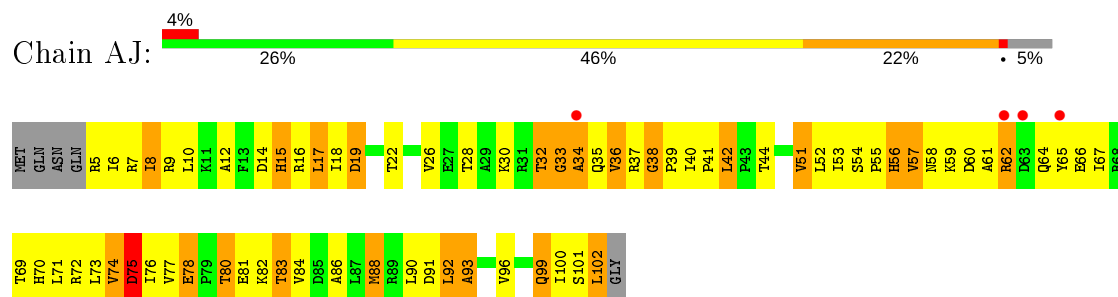
- Molecule 8: 30S ribosomal protein S9



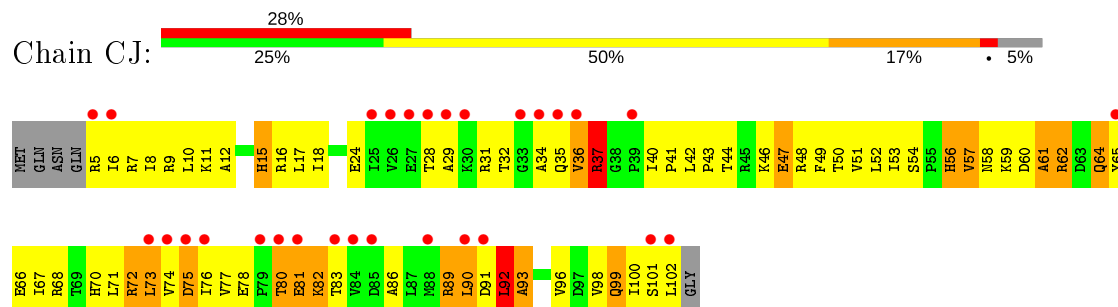
• Molecule 8: 30S ribosomal protein S9



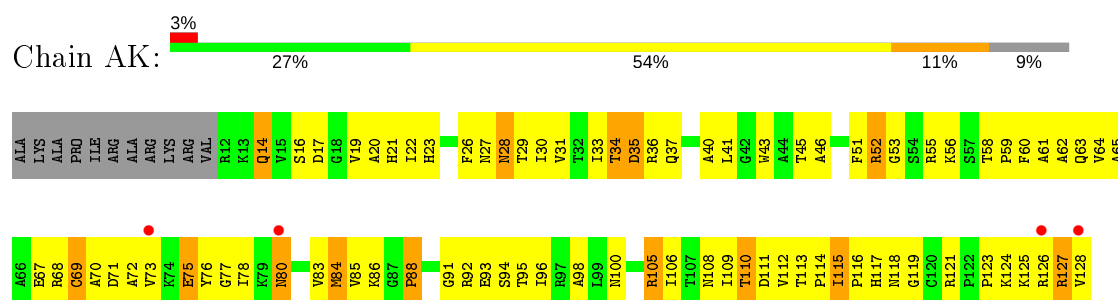
• Molecule 9: 30S ribosomal protein S10



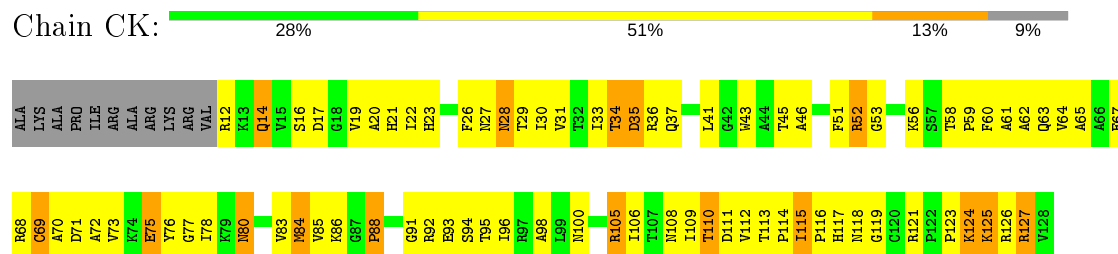
• Molecule 9: 30S ribosomal protein S10



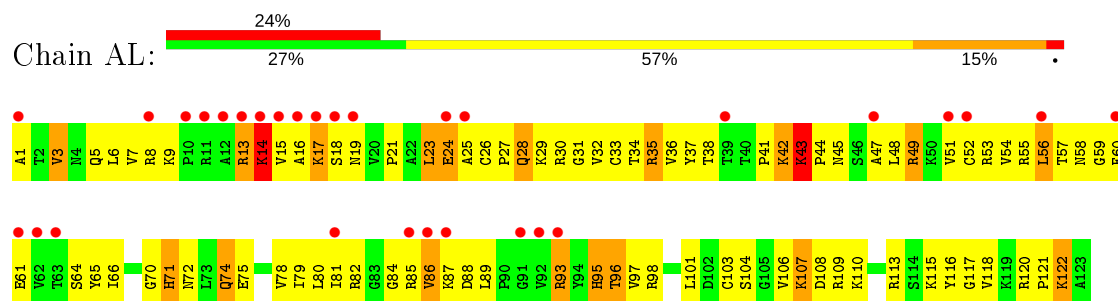
• Molecule 10: 30S ribosomal protein S11



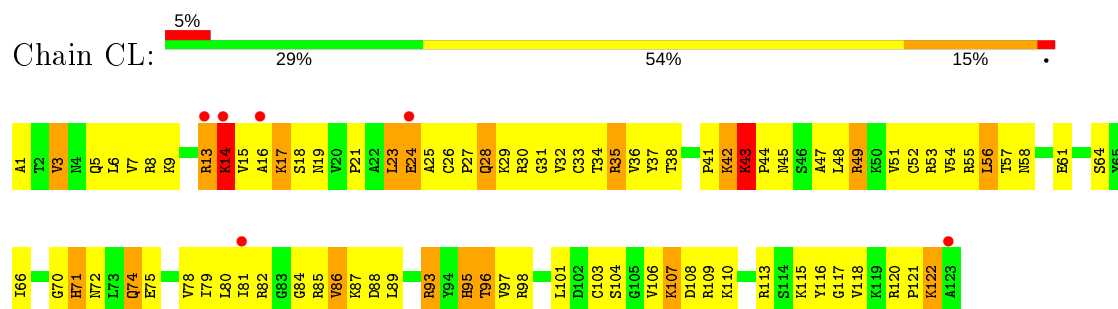
- Molecule 10: 30S ribosomal protein S11



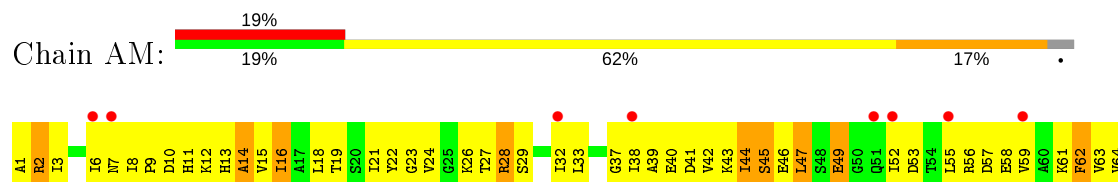
- Molecule 11: 30S ribosomal protein S12

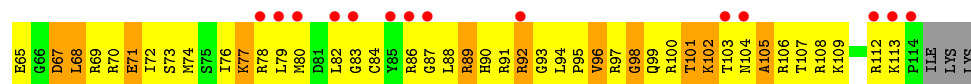


- Molecule 11: 30S ribosomal protein S12



- Molecule 12: 30S ribosomal protein S13

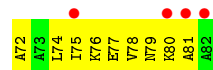
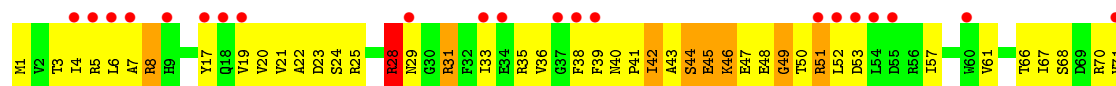




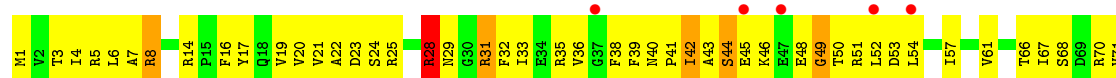
- Molecule 12: 30S ribosomal protein S13



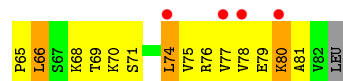
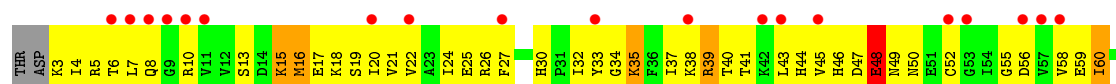
- Molecule 13: 30S ribosomal protein S16



- Molecule 13: 30S ribosomal protein S16

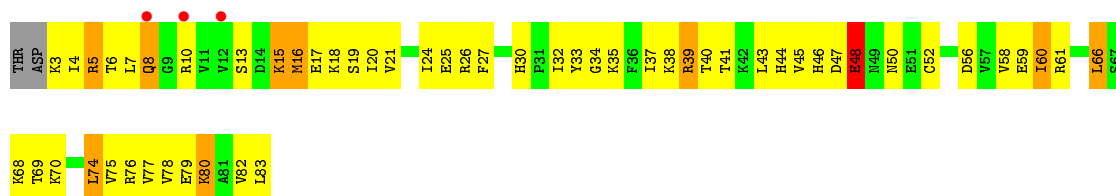


- Molecule 14: 30S ribosomal protein S17

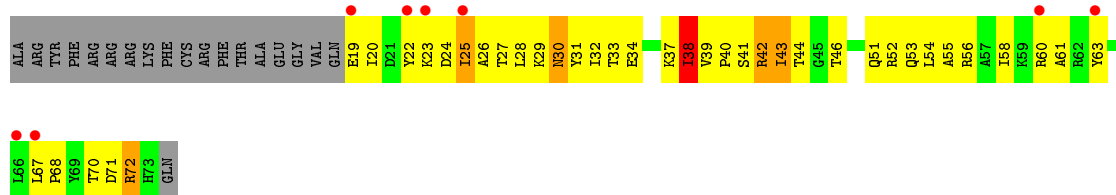
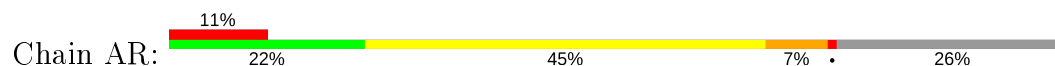


- Molecule 14: 30S ribosomal protein S17

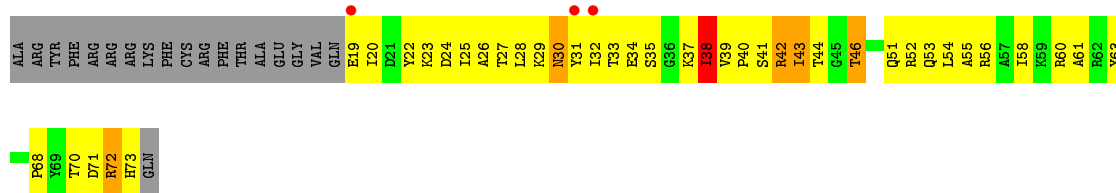
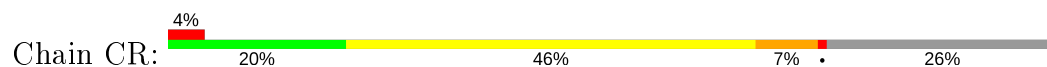




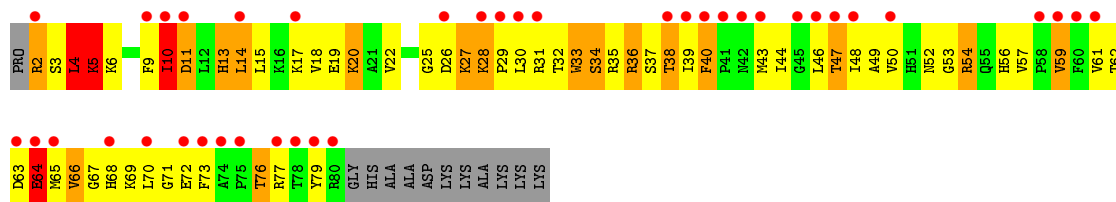
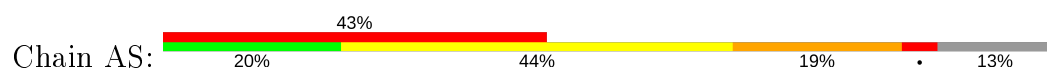
- Molecule 15: 30S ribosomal protein S18



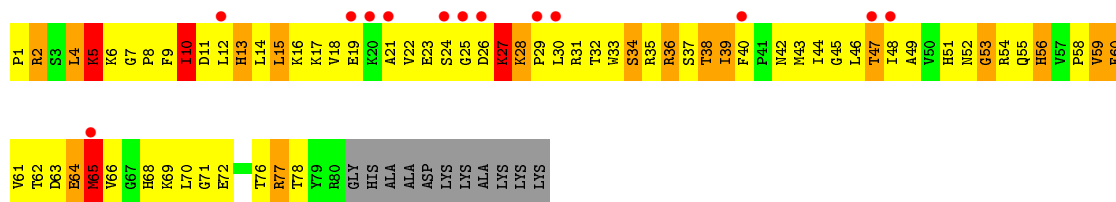
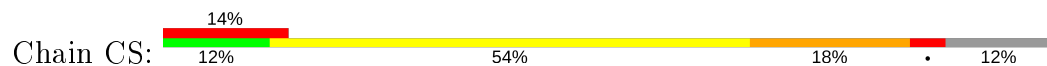
- Molecule 15: 30S ribosomal protein S18



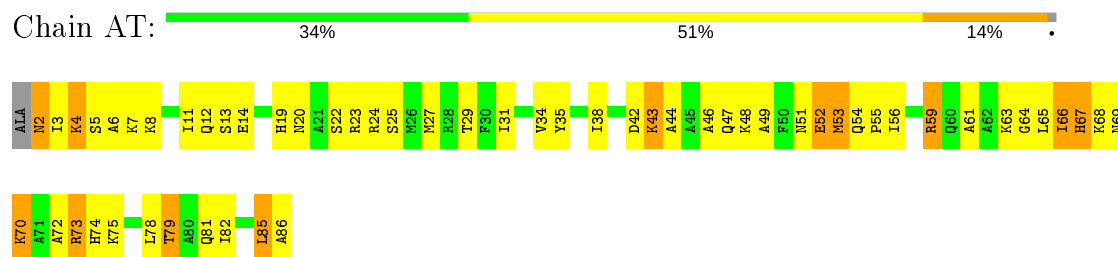
- Molecule 16: 30S ribosomal protein S19



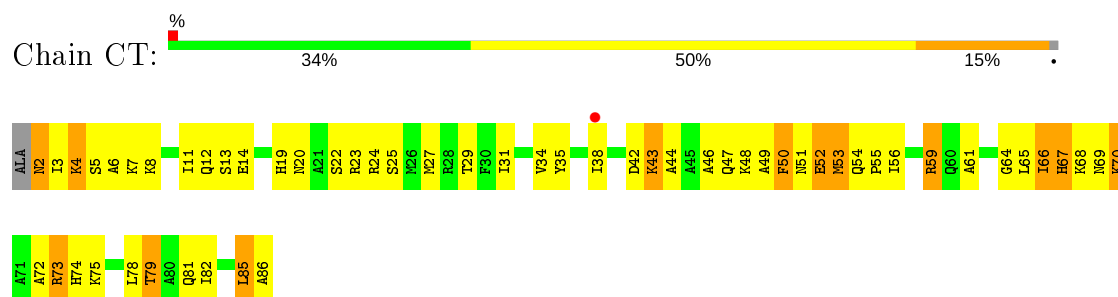
- Molecule 16: 30S ribosomal protein S19



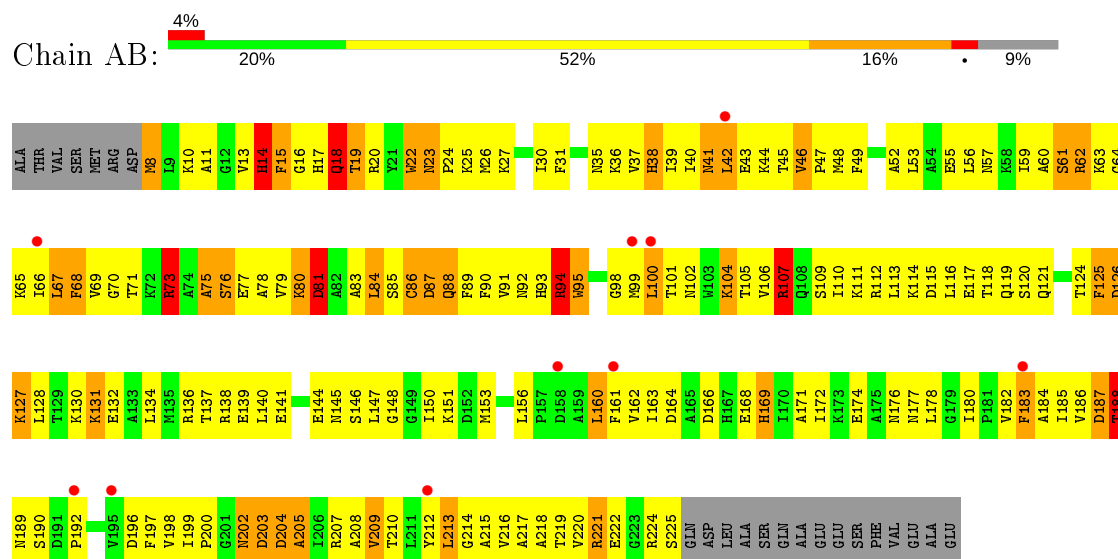
- Molecule 17: 30S ribosomal protein S20



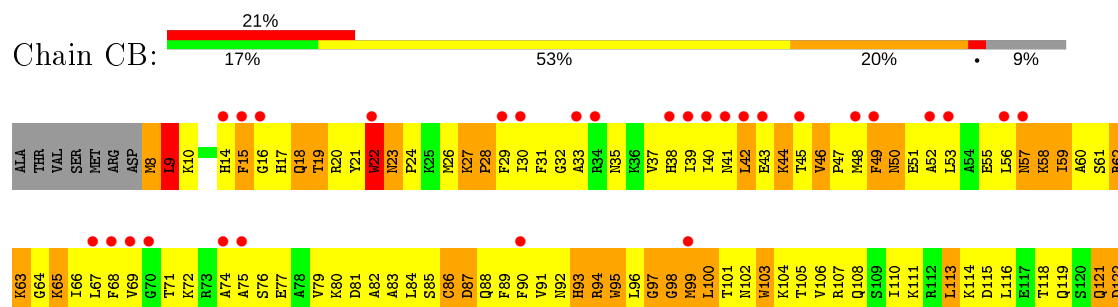
- Molecule 17: 30S ribosomal protein S20

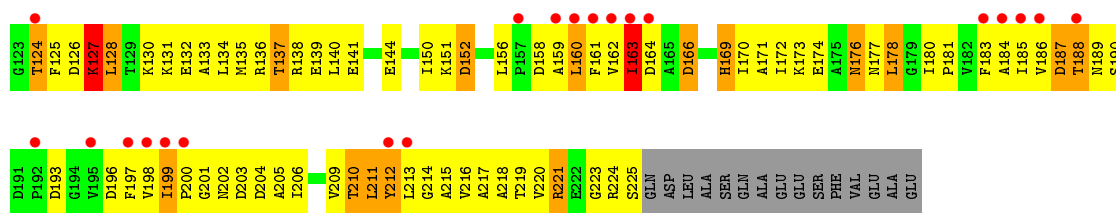


- Molecule 18: 30S ribosomal protein S2

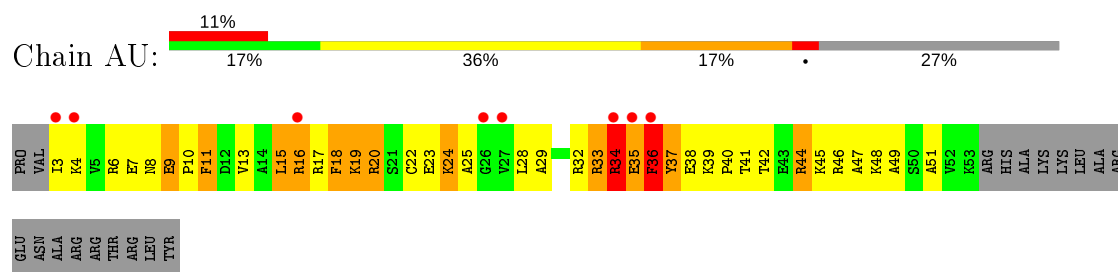


- Molecule 18: 30S ribosomal protein S2

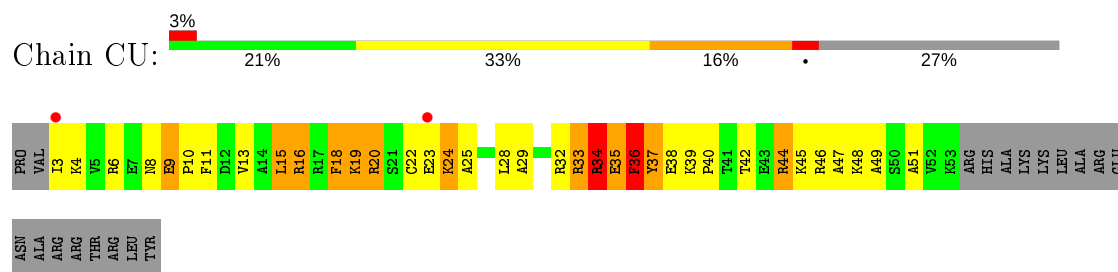




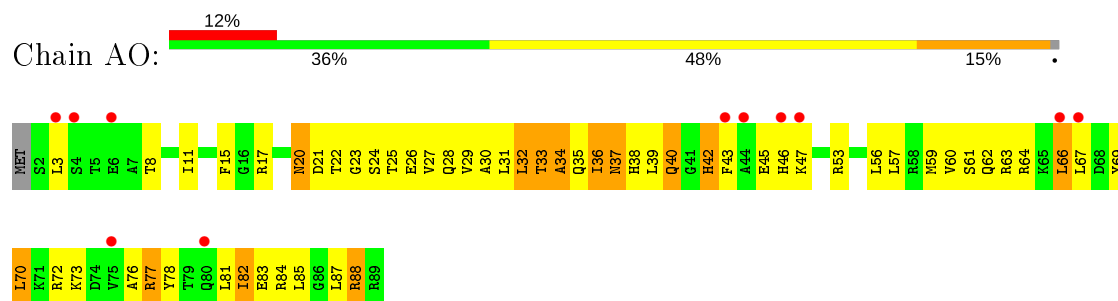
• Molecule 19: 30S ribosomal protein S21



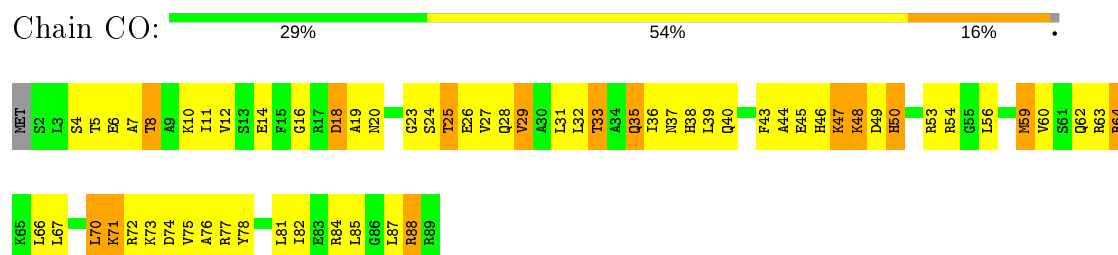
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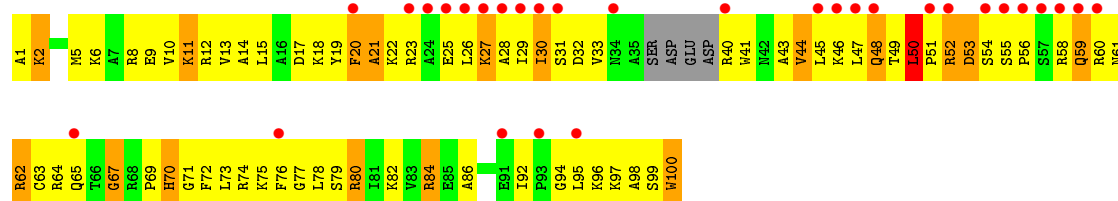
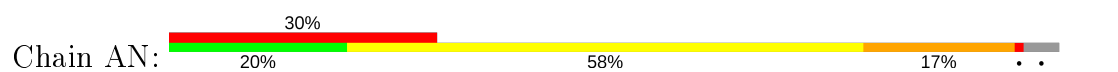
• Molecule 20: 30S ribosomal protein S15



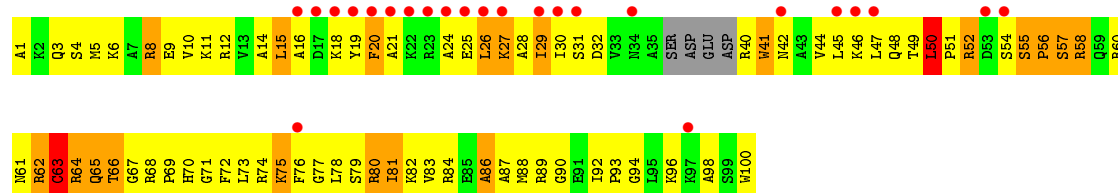
• Molecule 20: 30S ribosomal protein S15



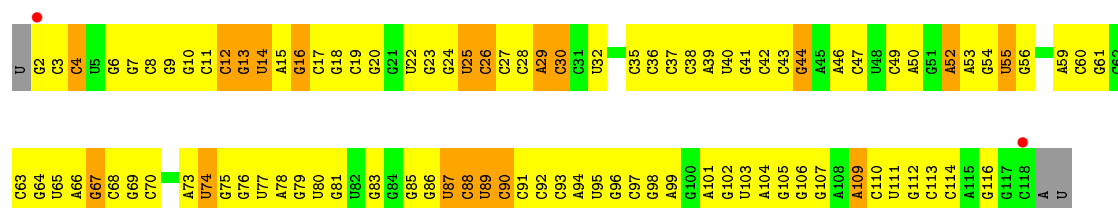
• Molecule 21: 30S ribosomal protein S14



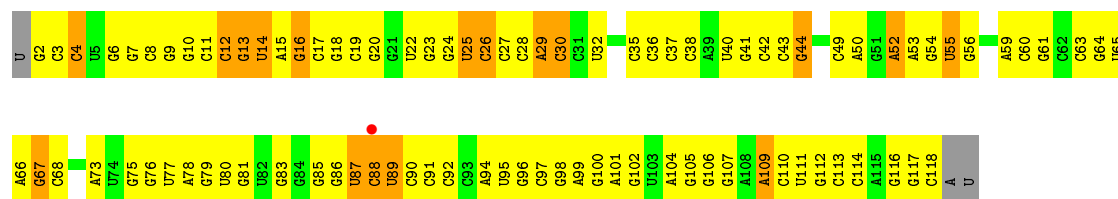
- Molecule 21: 30S ribosomal protein S14



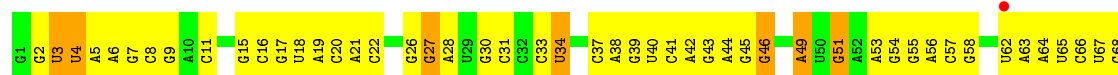
- Molecule 22: 5S rRNA



- Molecule 22: 5S rRNA



- Molecule 23: 23S rRNA

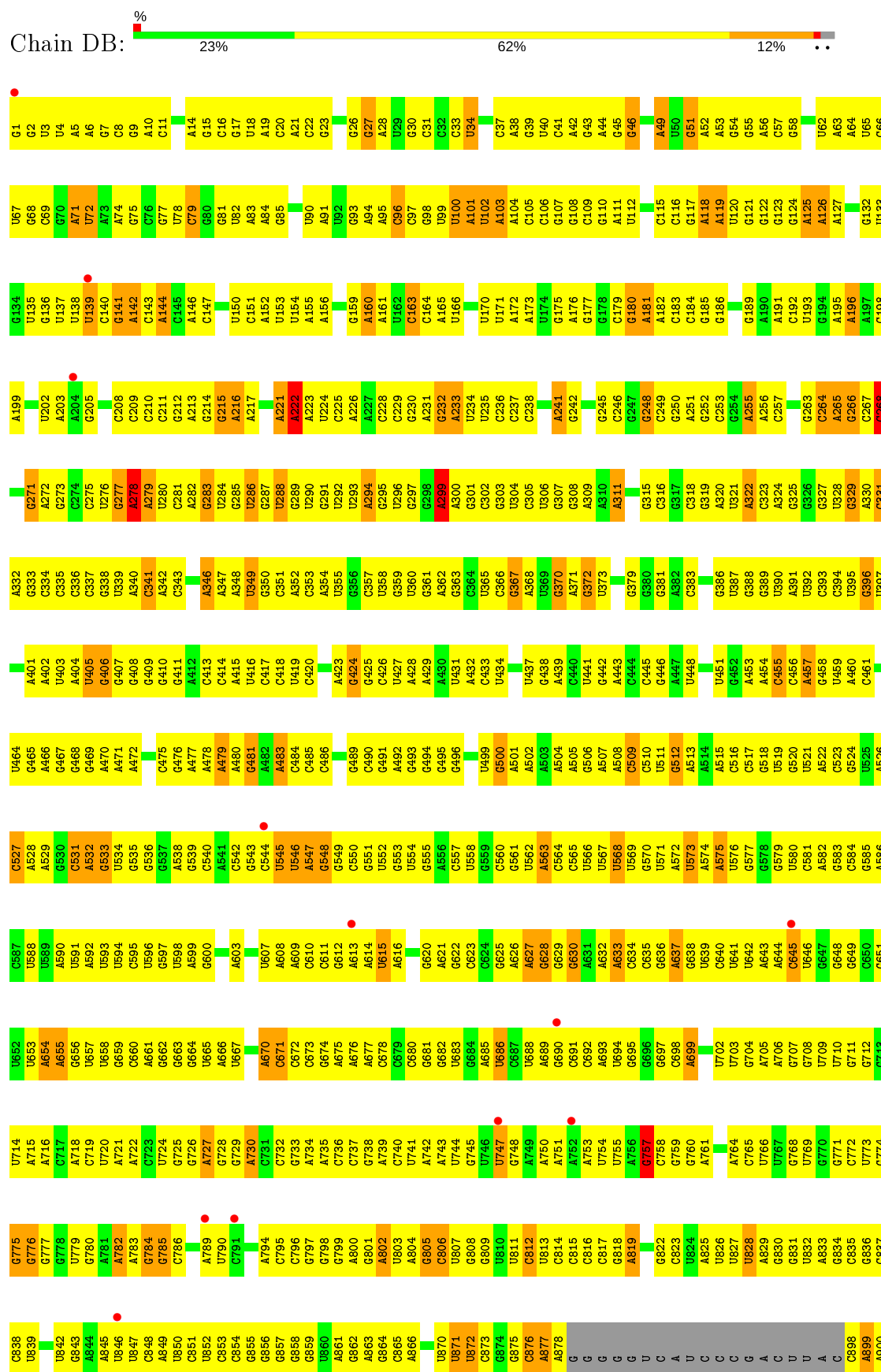


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G969	G907	G843	G778	A715	U852	A590	A828	G465	U403	C394	C268	U200	G136	G70
U970	G908	A844	U779	A716	U853	U591	A529	A466	A404	C395		C201	U137	A71
G971	G970	A845	G780	C717	A854	A592	G530	G467	U405	C396	G271	U202	U138	U72
A972	A909	U846	A781	A718	A855	U593	C531	G468	G406	G337	G272	A203	U139	A73
A973	A910	U847	A782	C719	G856	U594	A532	G469	G407	G338	G273	A204	C140	A74
G974	A911	C848	A783	U720	U857	C595	G533	A470	G408	U339	C274	G205	G141	G75
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	U913	U850	G785	A722	G859	G597		A472	G410	C341	U276	C208	C143	G77
	C786			G723	C851	U598	A538		A411	A342	G277	C209	A144	U78
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C982	A918	G855	A727	G664	G665	A603	C543	A478	A415	A346	C281	A213	U148	U82
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C986	C922	G859				U607	U547	A483	U419	C351	G287	A217	A152	
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A990	G926	A863	C737	G873	C550	C611	C550		A423	U355	U292	A223	G93	G93
C991	A927	G864	G738	G874	G551	G612	G551		G424		U293	U224	A94	A94
C992	C865	A865	A675	G875	U552	A613	U552	G489	G425	G359	G295	C225	A95	A95
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C1001	C937	A877	G748	G884	G561	G623	G561	U499	U434	G367	G303	U234	U170	A104
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C1012	A945	U	U755	C691	U569	A631	U569	A507	A443	G380	A311	U243	G177	U113
C1013	C946	C	A756	C692	A632	G632	G570	A508	C444	G381	G315	G244	G180	C116
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G1016	G949	C	G760	C698	U697	G636	U573	U511	A447	C386	G318	G247	C183	A119
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A1020	U955	A	C765	U702	U703	U639	G577	A515	G452	A391	A322	G252	G188	G123
C1021	G956	U	U766	G703	G704	C640	G578	C516	A453	A392	G323	G253	G189	G124
G1022	C957	C	U767	G704	G705	U641	U580	C517	A454	A393	C324	G254	A190	A126
U1023	U958	U	G768	A705	A706	A642	U581	C518	C455	U392	A324	A255	A191	A127
G1024	A959	A	U769	A706	A707	A643	A582	U519	C456	C393	G325	A256	C192	C128
G1025	C960	C	G770	A707	A708	A644	U583	G520	A457	C394	G326	A257	G193	G129
A1026	U960	C	G771	G708	G709	C645	U521	G458	U459	U395	U327	C257	U193	C130
G1027	C961	C	G772	G709	U709	U646	C584	A522	U459	G396	U328		G194	
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C1029	U963	C	G774	G711	G712	G648	A586	U524	C461	G400	A330	C264	A196	G132
U1030			G775	G712	G713	C850	C587	U525	C462		C331	A265	A197	U133
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- Molecule 23: 23S rRNA

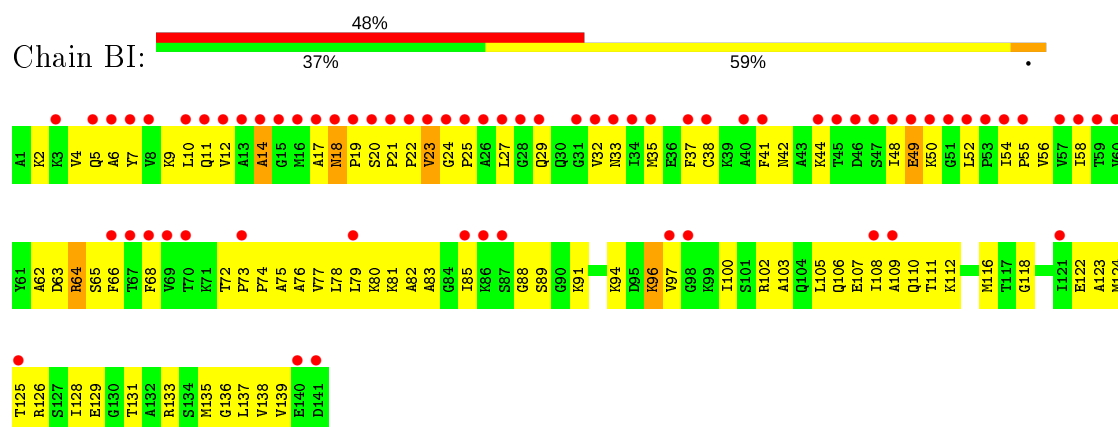


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C1795	C1795	G1666	U1598	G1474	C1414	G1350	A1284	U1219	C1150	U1083	G1011	C946
U1796	U1796	A1669	A1599	U1475	U1415	C1351	A1285	G1220	A1151	A1084	U1012	A947
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C1800	C1800	G1673	A1603	U1479	A1418	G1355	C1289	G1225	G1157	A1088	G1016	C951
A1801	A1801	U1674	U1603	G1480	U1419	U1356	G1290	G1226	C1158	A1089	G1017	G952
G1738	G1738	G1675	C1607	U1481	A1420	G1357	A1289	A1227	G1159	A1090	U1018	U955
A1802	A1802	C1676	A1608	U1482	G1421	U1359	G1291	G1228	C1161	G1091	U1019	U956
C1803	C1803	U1677	A1609	U1483	G1422	G1360	G1292	G1229	G1162	C1092	A1020	G957
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G1807	G1807	A1680	U1615	U1487	G1426	A1365	G1296	G1237	G1166	A1098	G1024	C961
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G1813	G1813	C1686	U1621	U1493								
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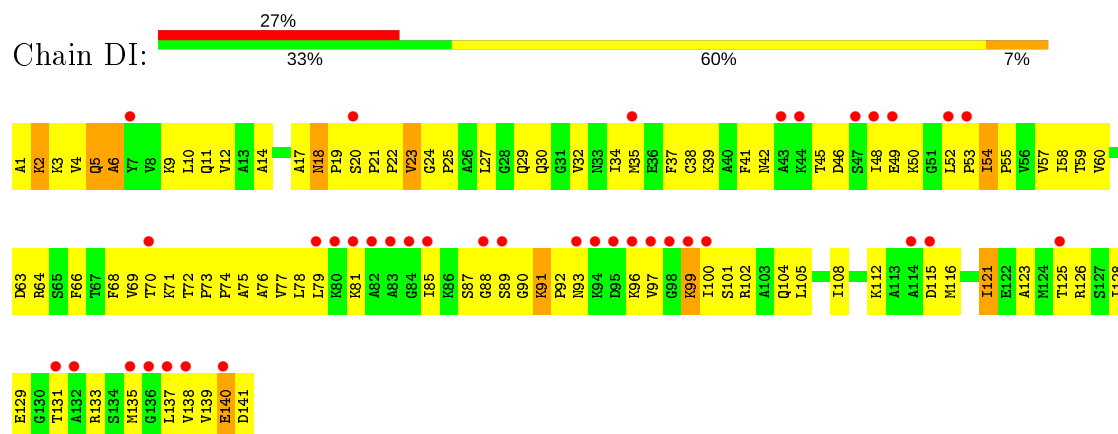
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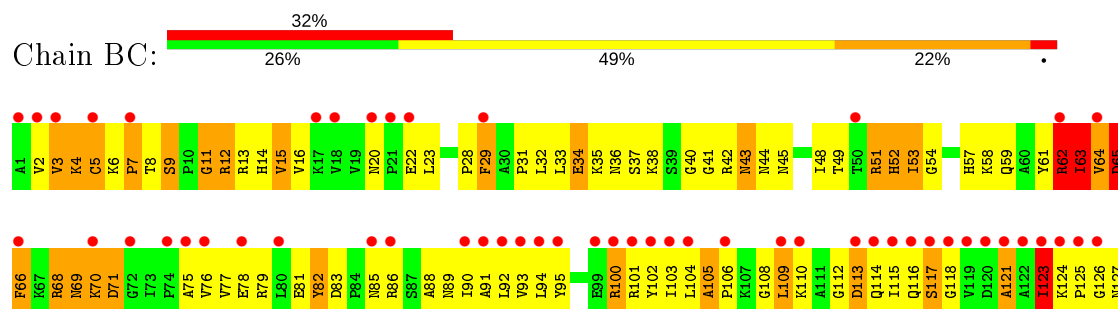
• Molecule 24: 50S ribosomal protein L11

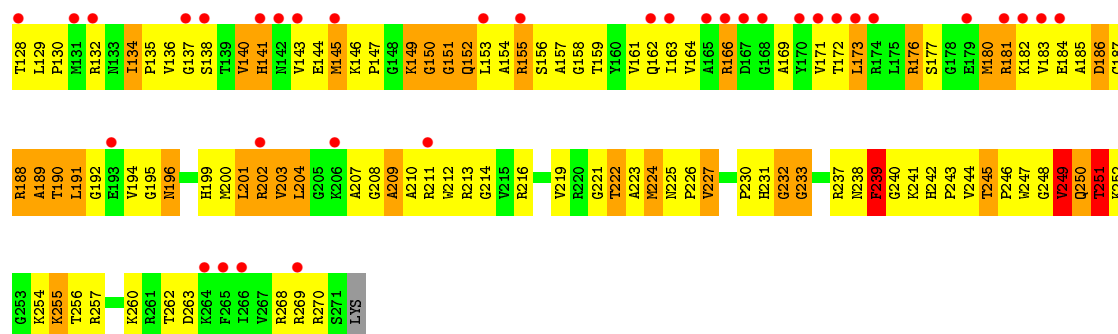


• Molecule 24: 50S ribosomal protein L11

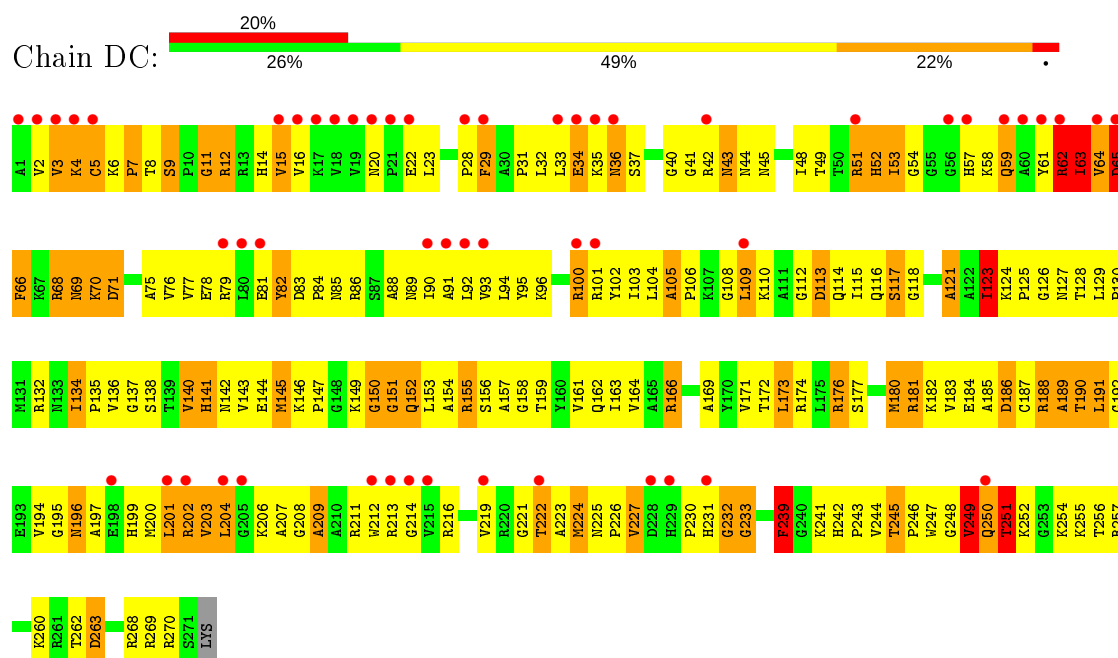


• Molecule 25: 50S ribosomal protein L2

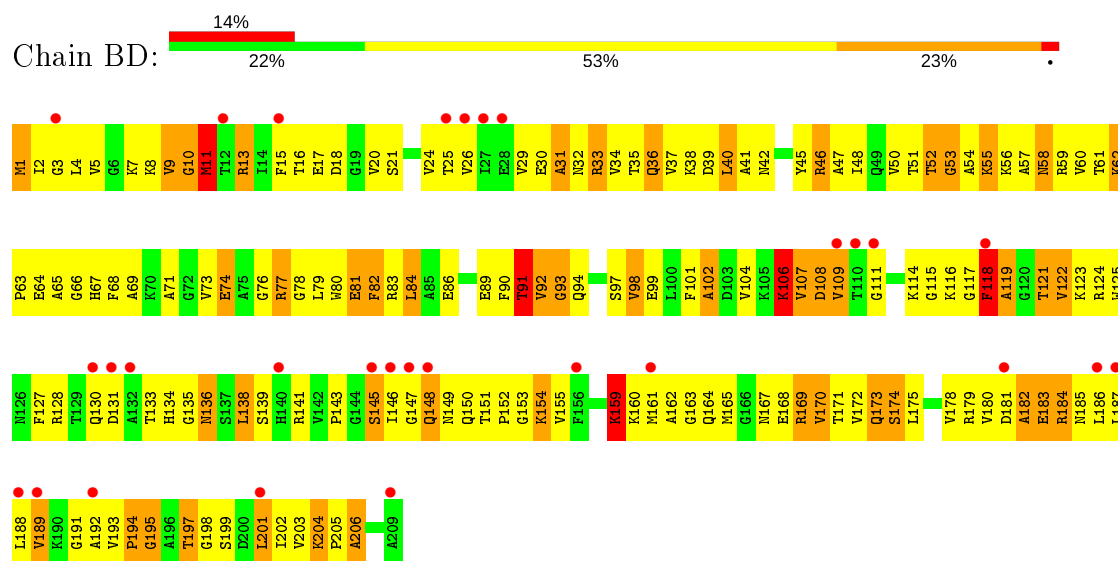




• Molecule 25: 50S ribosomal protein L2

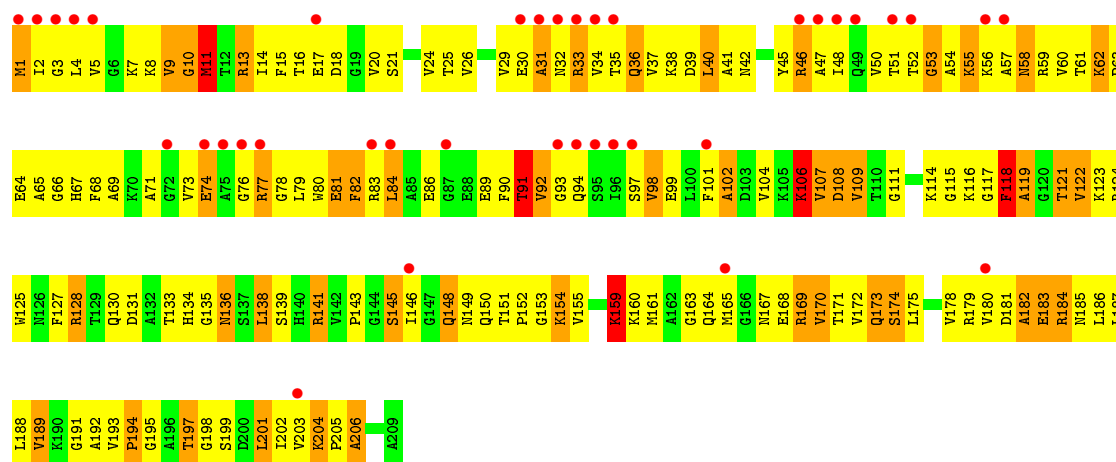


• Molecule 26: 50S ribosomal protein L3

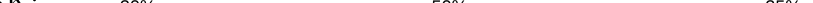


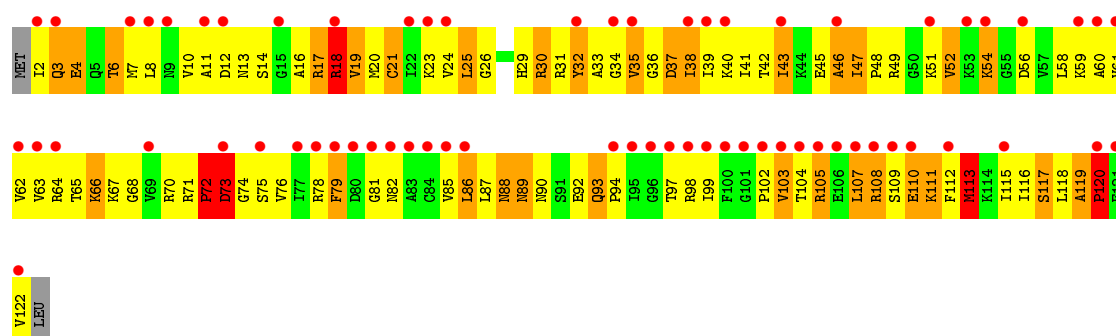
• Molecule 26: 50S ribosomal protein L3

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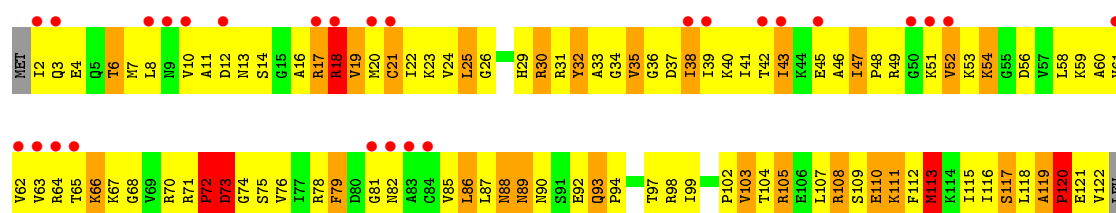
- Molecule 27: 50S ribosomal protein L14

Chain BK: 



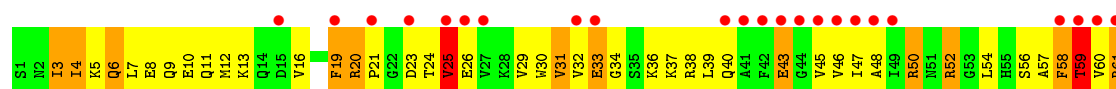
- Molecule 27: 50S ribosomal protein L14

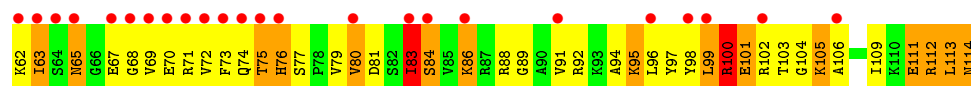
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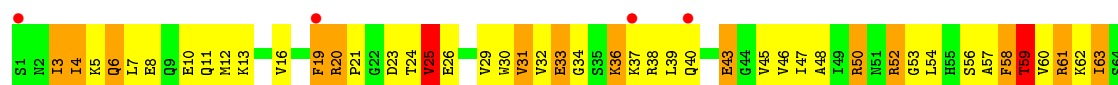
- Molecule 28: 50S ribosomal protein L19

Chain BP: 

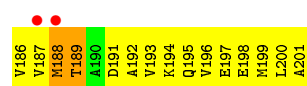
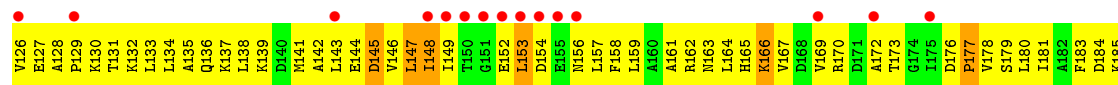
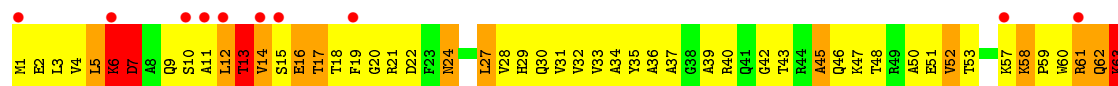




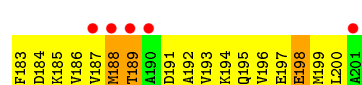
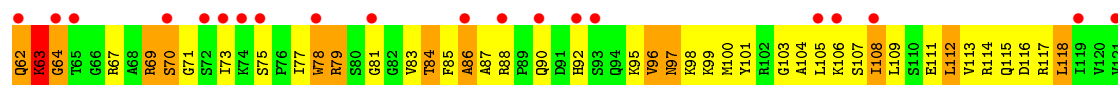
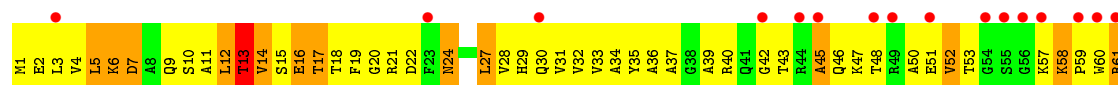
• Molecule 28: 50S ribosomal protein L19



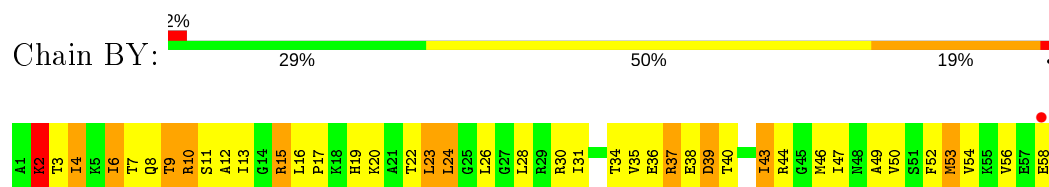
• Molecule 29: 50S ribosomal protein L4



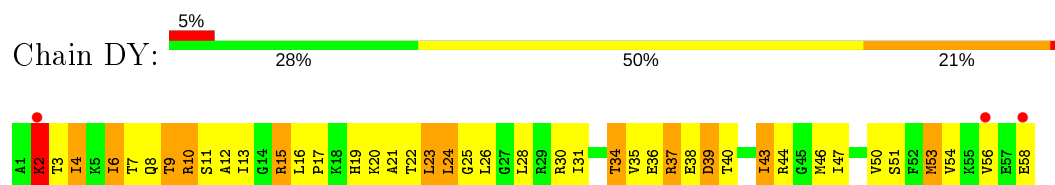
• Molecule 29: 50S ribosomal protein L4



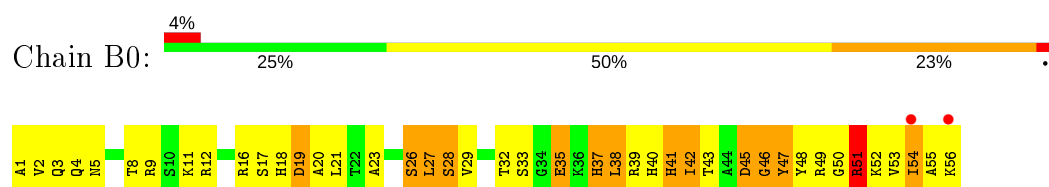
- Molecule 30: 50S ribosomal protein L30



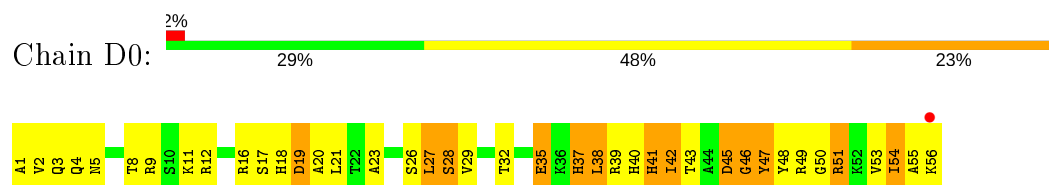
- Molecule 30: 50S ribosomal protein L30



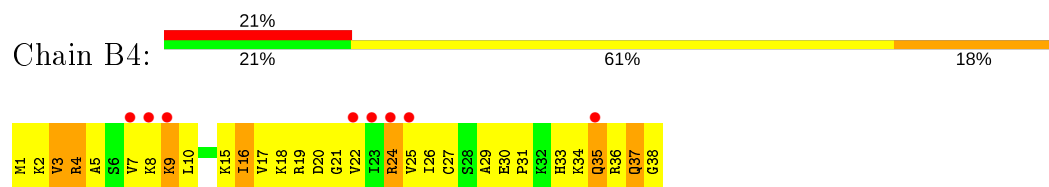
- Molecule 31: 50S ribosomal protein L32



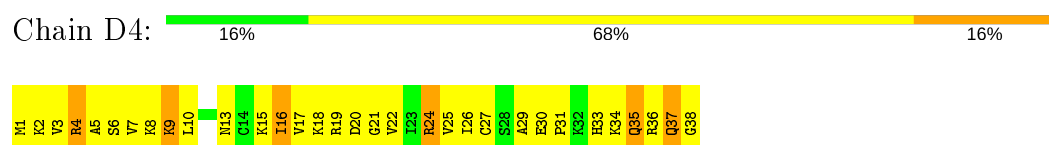
- Molecule 31: 50S ribosomal protein L32



- Molecule 32: 50S ribosomal protein L36



- Molecule 32: 50S ribosomal protein L36

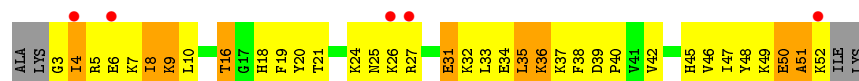


- Molecule 33: 50S ribosomal protein L33

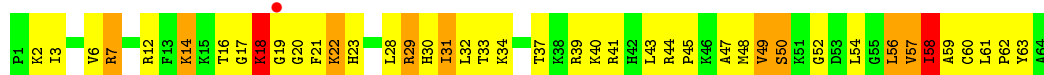




- Molecule 33: 50S ribosomal protein L33



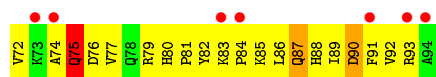
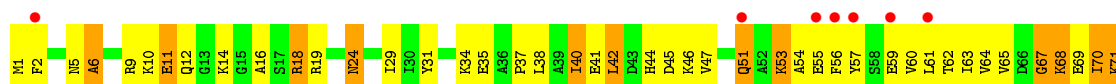
- Molecule 34: 50S ribosomal protein L35



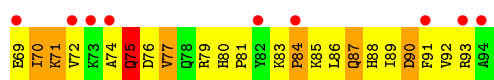
- Molecule 34: 50S ribosomal protein L35



- Molecule 35: 50S ribosomal protein L25

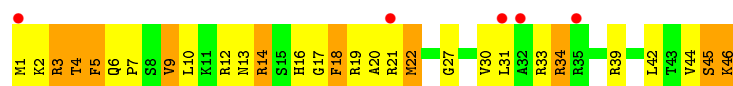


- Molecule 35: 50S ribosomal protein L25

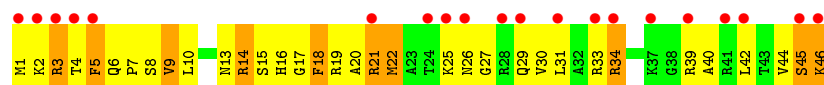


- Molecule 36: 50S ribosomal protein L34

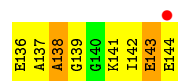




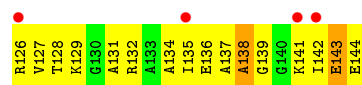
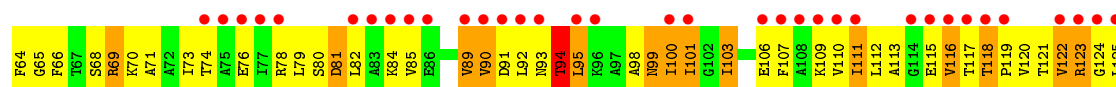
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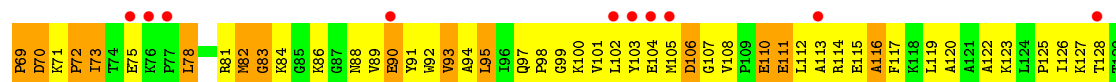
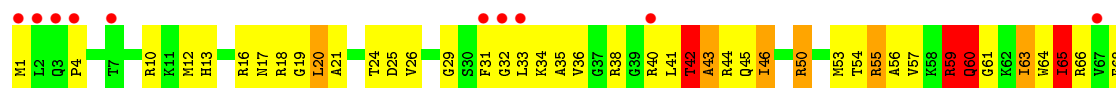
• Molecule 37: 50S ribosomal protein L15

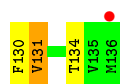


• Molecule 37: 50S ribosomal protein L15

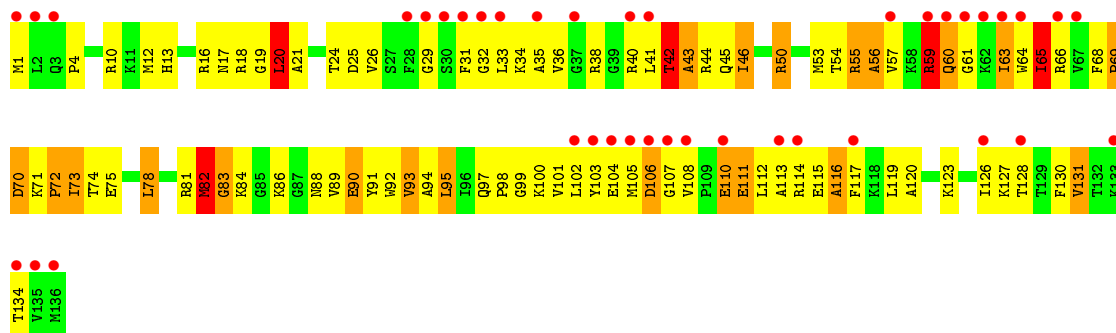


• Molecule 38: 50S ribosomal protein L16

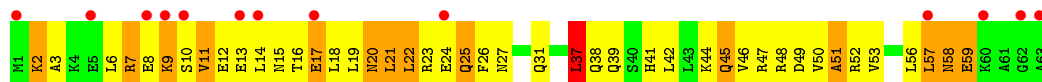




• Molecule 38: 50S ribosomal protein L16



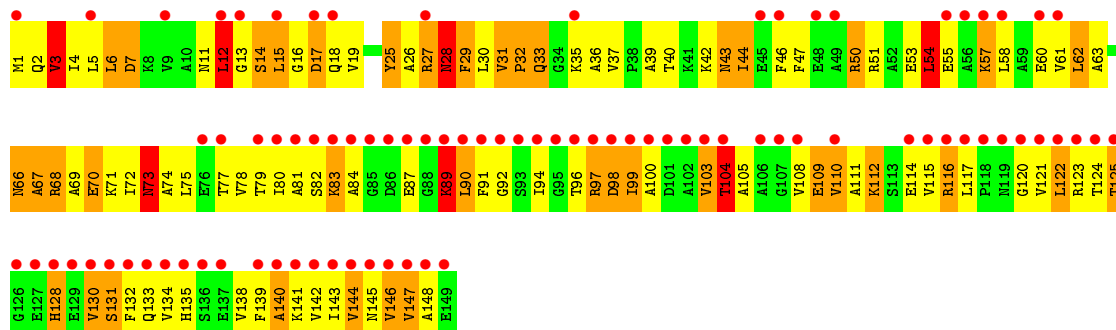
• Molecule 39: 50S ribosomal protein L29



• Molecule 39: 50S ribosomal protein L29

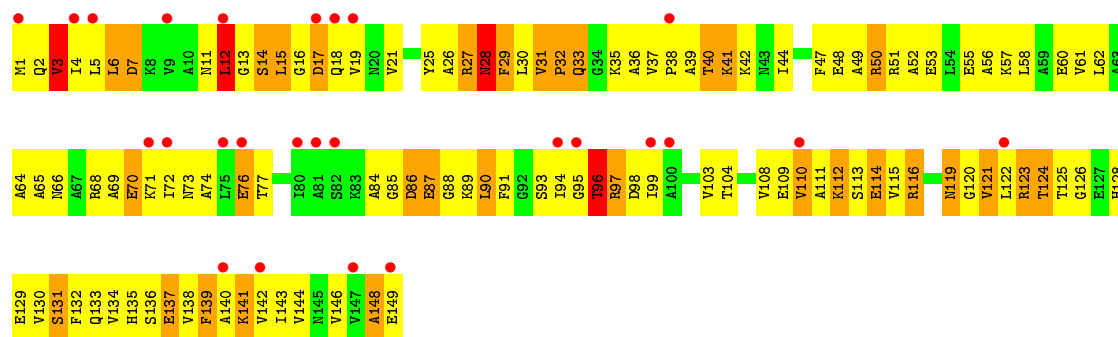


• Molecule 40: 50S ribosomal protein L9

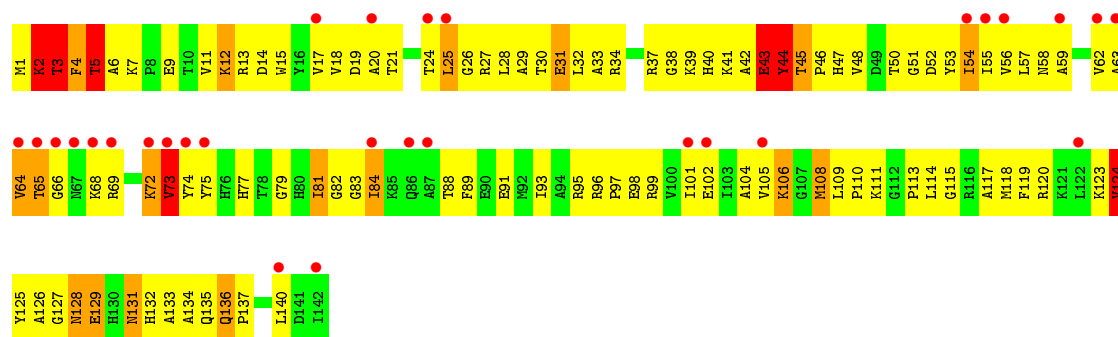


• Molecule 40: 50S ribosomal protein L9

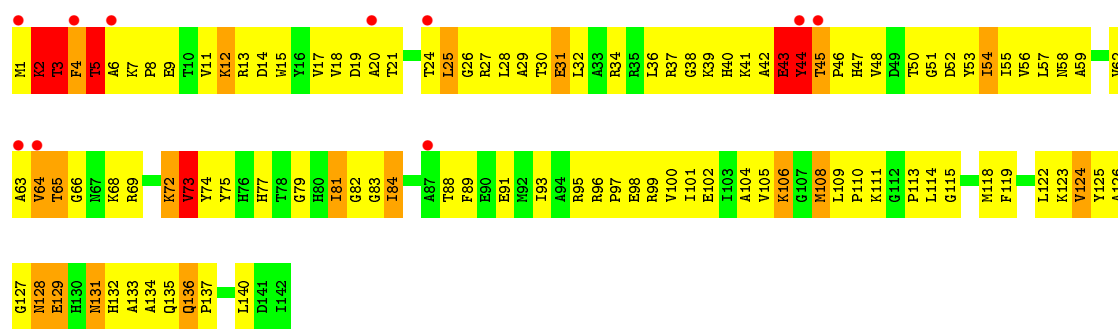




● Molecule 41: 50S ribosomal protein L13



● Molecule 41: 50S ribosomal protein L13



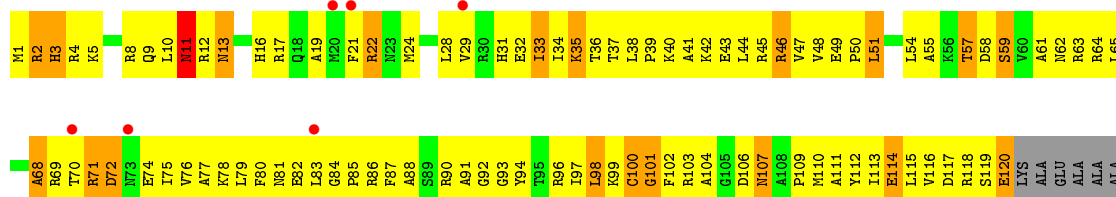
● Molecule 42: 50S ribosomal protein L17



GLU

- Molecule 42: 50S ribosomal protein L17

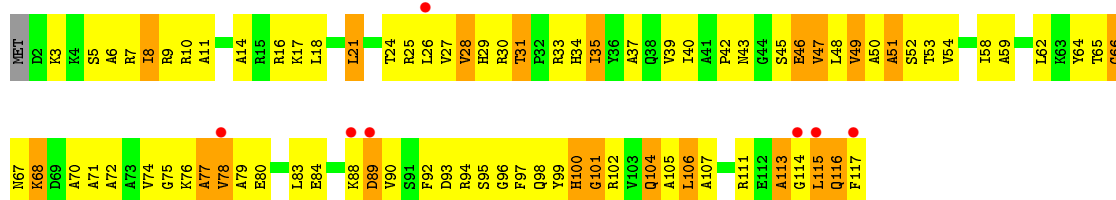
Chain DN: 5% 17% 61% 15% 6%



GLU

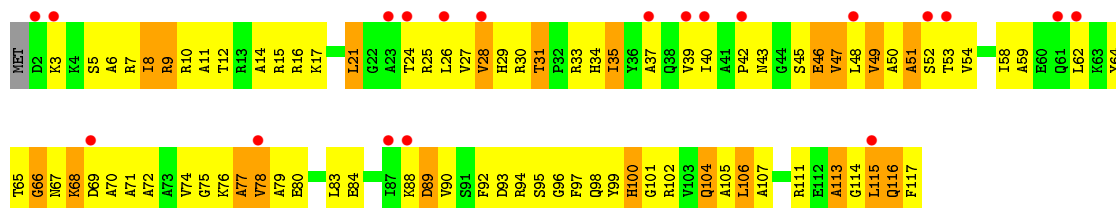
- Molecule 43: 50S ribosomal protein L18

Chain BO: 6% 28% 53% 18%



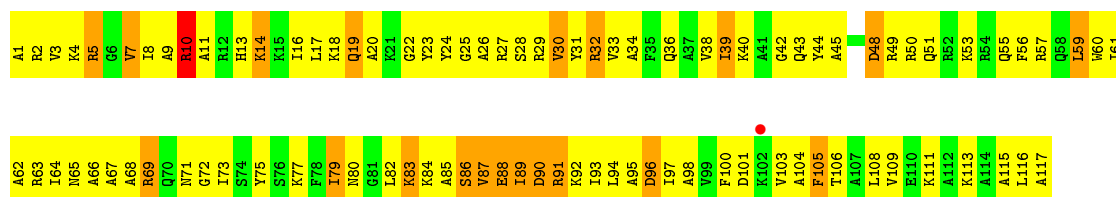
- Molecule 43: 50S ribosomal protein L18

Chain DO: 17% 26% 55% 18%

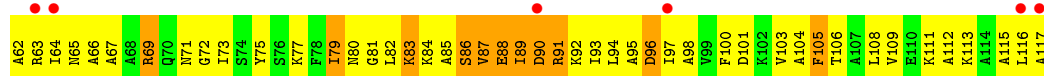
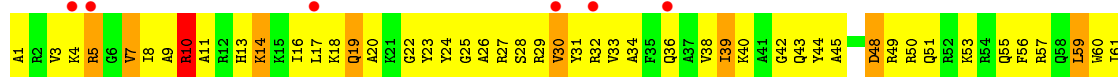


- Molecule 44: 50S ribosomal protein L20

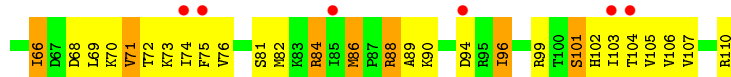
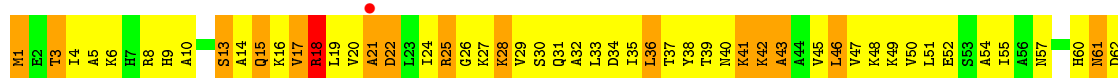
Chain BQ: 20% 62% 17%



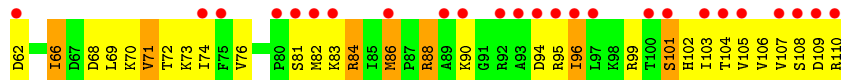
- Molecule 44: 50S ribosomal protein L20



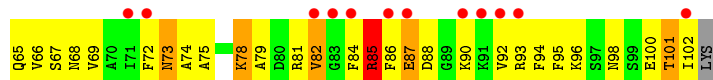
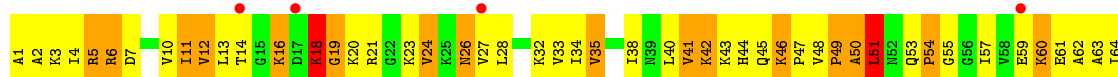
• Molecule 45: 50S ribosomal protein L22



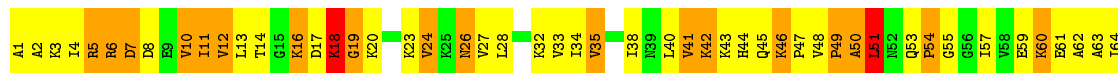
• Molecule 45: 50S ribosomal protein L22



• Molecule 46: 50S ribosomal protein L24

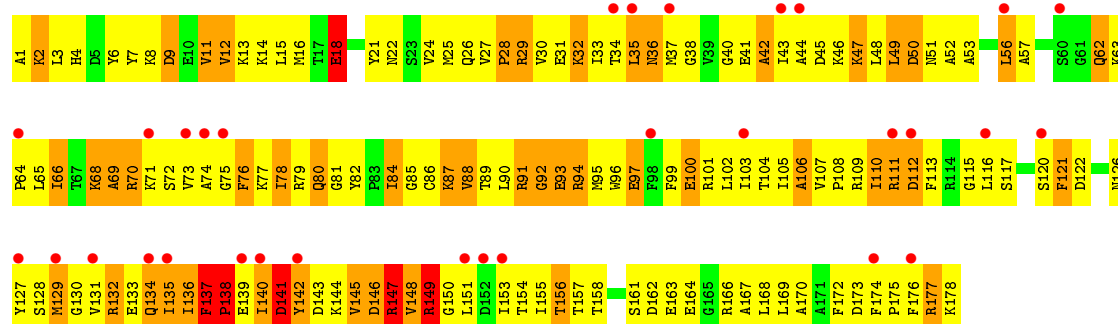
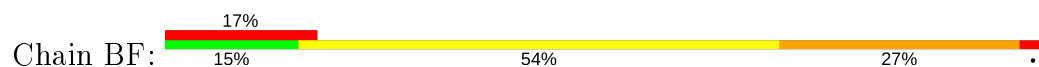


• Molecule 46: 50S ribosomal protein L24

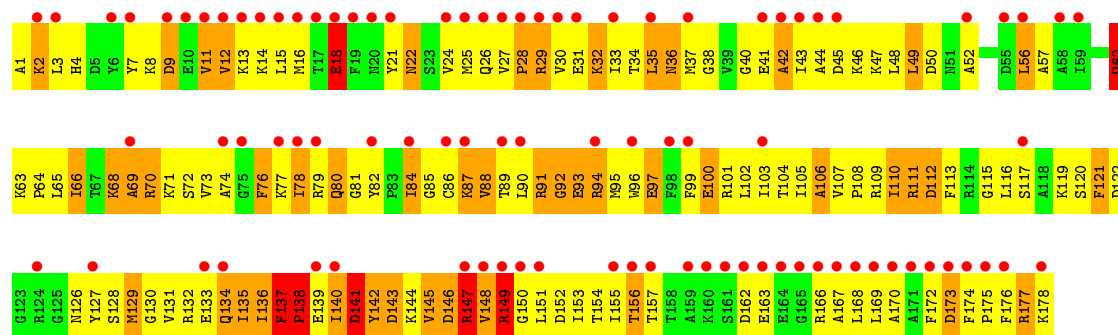
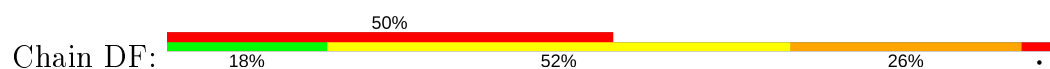




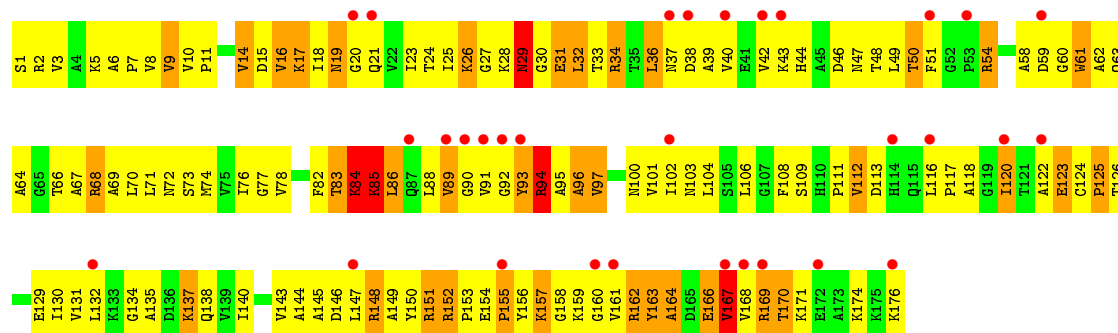
• Molecule 47: 50S ribosomal protein L5



• Molecule 47: 50S ribosomal protein L5

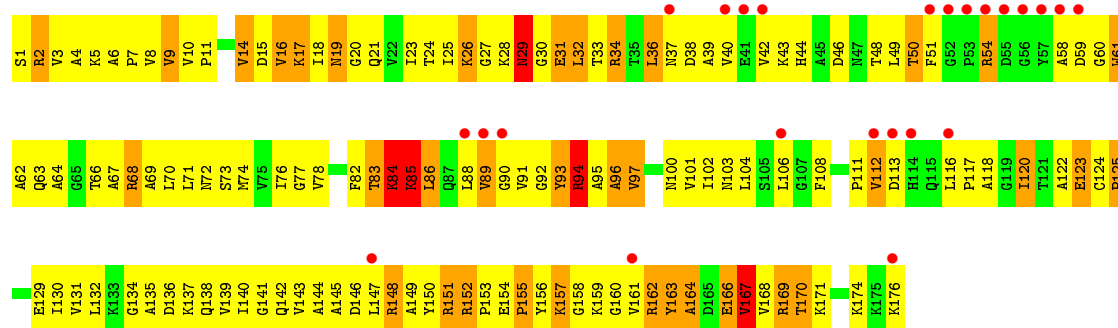


• Molecule 48: 50S ribosomal protein L6

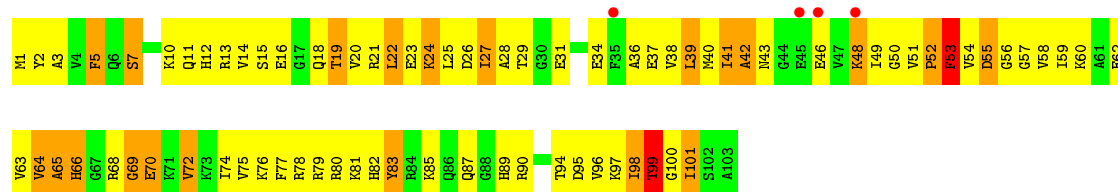


• Molecule 48: 50S ribosomal protein L6

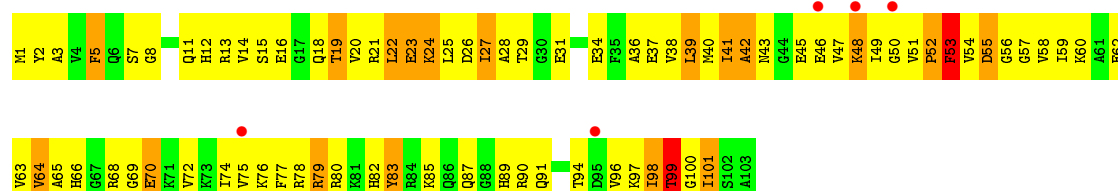




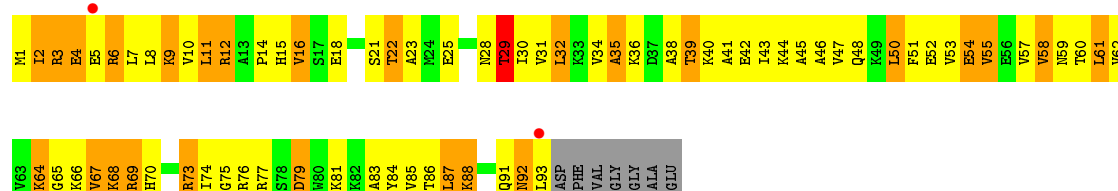
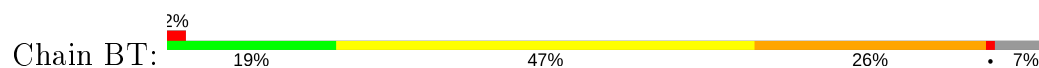
• Molecule 49: 50S ribosomal protein L21



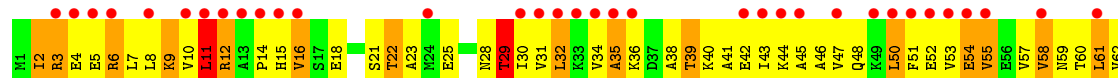
• Molecule 49: 50S ribosomal protein L21



• Molecule 50: 50S ribosomal protein L23

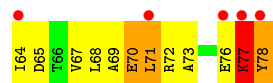
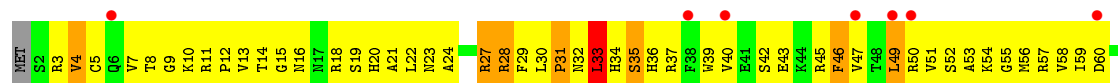


• Molecule 50: 50S ribosomal protein L23

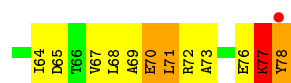
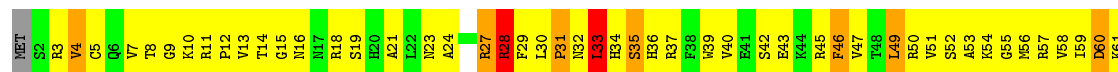




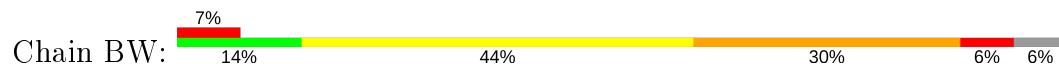
- Molecule 51: 50S ribosomal protein L28



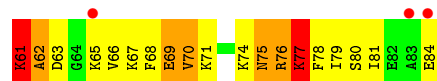
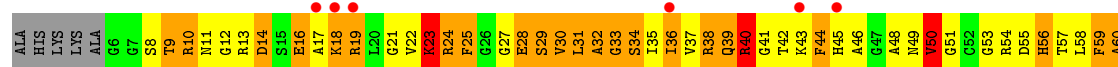
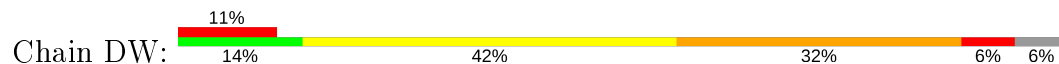
- Molecule 51: 50S ribosomal protein L28



- Molecule 52: 50S ribosomal protein L27



- Molecule 52: 50S ribosomal protein L27



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	208.85Å 379.20Å 739.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.00 – 3.93 182.94 – 3.94	Depositor EDS
% Data completeness (in resolution range)	(Not available) (70.00-3.93) 75.9 (182.94-3.94)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 3.89Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.258 , 0.311 0.224 , 0.271	Depositor DCC
R_{free} test set	19247 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	149.7	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.20 , 71.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	284033	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.50% 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: SCM, 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	0.27	2/36762 (0.0%)	0.76	12/57350 (0.0%)
1	CA	0.31	1/36762 (0.0%)	0.77	11/57350 (0.0%)
2	AC	0.23	0/1651	0.45	0/2225
2	CC	0.23	0/1651	0.46	0/2225
3	AD	0.23	0/1665	0.44	0/2227
3	CD	0.23	0/1665	0.44	0/2227
4	AE	0.23	0/1118	0.45	0/1504
4	CE	0.23	0/1118	0.45	0/1504
5	AF	0.24	0/835	0.45	0/1128
5	CF	0.24	0/835	0.45	0/1128
6	AG	0.23	0/1187	0.45	0/1591
6	CG	0.23	0/1211	0.45	0/1624
7	AH	0.23	0/989	0.44	0/1326
7	CH	0.23	0/989	0.44	0/1326
8	AI	0.24	0/1034	0.44	0/1375
8	CI	0.24	0/1034	0.45	0/1375
9	AJ	0.22	0/796	0.48	0/1077
9	CJ	0.22	0/796	0.48	0/1077
10	AK	0.24	0/893	0.44	0/1205
10	CK	0.24	0/893	0.44	0/1205
11	AL	0.22	0/969	0.48	0/1300
11	CL	0.22	0/969	0.48	0/1300
12	AM	0.21	0/892	0.45	0/1193
12	CM	0.21	0/884	0.45	0/1181
13	AP	0.25	0/659	0.45	0/884
13	CP	0.25	0/648	0.44	0/870
14	AQ	0.23	0/657	0.46	0/881
14	CQ	0.24	0/666	0.46	0/892
15	AR	0.23	0/462	0.46	0/621
15	CR	0.23	0/462	0.46	0/621
16	AS	0.25	0/652	0.46	0/877
16	CS	0.25	0/660	0.49	0/888

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AT	0.23	0/671	0.40	0/888
17	CT	0.23	0/671	0.40	0/888
18	AB	0.25	0/1735	0.45	0/2338
18	CB	0.25	0/1735	0.45	0/2338
19	AU	0.26	0/430	0.46	0/570
19	CU	0.25	0/430	0.46	0/570
20	AO	0.22	0/722	0.45	0/964
20	CO	0.23	0/722	0.44	0/964
21	AN	0.24	0/785	0.44	0/1043
21	CN	0.24	0/785	0.46	0/1043
22	BA	0.23	0/2803	0.74	1/4371 (0.0%)
22	DA	0.24	0/2803	0.75	1/4371 (0.0%)
23	BB	0.28	5/68314 (0.0%)	0.77	33/106569 (0.0%)
23	DB	0.28	5/68314 (0.0%)	0.77	33/106569 (0.0%)
24	BI	0.24	0/1046	0.46	0/1410
24	DI	0.25	0/1046	0.47	0/1410
25	BC	0.22	0/2121	0.47	0/2852
25	DC	0.22	0/2121	0.47	0/2852
26	BD	0.24	0/1586	0.48	0/2134
26	DD	0.24	0/1586	0.48	0/2134
27	BK	0.24	0/939	0.53	0/1258
27	DK	0.24	0/939	0.53	0/1258
28	BP	0.24	0/929	0.49	0/1242
28	DP	0.24	0/929	0.49	0/1242
29	BE	0.24	0/1571	0.48	0/2113
29	DE	0.24	0/1571	0.48	0/2113
30	BY	0.24	0/453	0.49	0/605
30	DY	0.23	0/453	0.49	0/605
31	B0	0.23	0/450	0.51	0/599
31	D0	0.23	0/450	0.51	0/599
32	B4	0.22	0/303	0.49	0/397
32	D4	0.23	0/303	0.49	0/397
33	B1	0.27	0/416	0.48	0/554
33	D1	0.27	0/416	0.48	0/554
34	B3	0.24	0/513	0.46	0/676
34	D3	0.24	0/513	0.46	0/676
35	BV	0.25	0/766	0.43	0/1025
35	DV	0.25	0/766	0.43	0/1025
36	B2	0.26	0/380	0.47	0/498
36	D2	0.26	0/380	0.47	0/498
37	BL	0.24	0/1054	0.48	0/1403
37	DL	0.24	0/1054	0.48	0/1403
38	BM	0.25	0/1093	0.47	0/1460

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	DM	0.25	0/1093	0.47	0/1460
39	BX	0.24	0/510	0.50	0/677
39	DX	0.23	0/510	0.50	0/677
40	BH	0.25	0/1122	0.47	0/1515
40	DH	0.25	0/1122	0.47	0/1515
41	BJ	0.23	0/1152	0.47	0/1551
41	DJ	0.23	0/1152	0.47	0/1551
42	BN	0.24	0/973	0.49	0/1301
42	DN	0.24	0/973	0.49	0/1301
43	BO	0.23	0/902	0.47	0/1209
43	DO	0.23	0/902	0.47	0/1209
44	BQ	0.25	0/960	0.47	0/1278
44	DQ	0.25	0/960	0.47	0/1278
45	BS	0.22	0/864	0.50	0/1156
45	DS	0.22	0/864	0.50	0/1156
46	BU	0.25	0/787	0.45	0/1051
46	DU	0.25	0/787	0.45	0/1051
47	BF	0.26	0/1444	0.49	0/1937
47	DF	0.26	0/1444	0.49	0/1937
48	BG	0.23	0/1343	0.47	0/1816
48	DG	0.23	0/1343	0.47	0/1816
49	BR	0.26	0/829	0.48	0/1107
49	DR	0.25	0/829	0.48	0/1107
50	BT	0.23	0/744	0.51	0/994
50	DT	0.23	0/744	0.51	0/994
51	BZ	0.25	0/635	0.49	0/848
51	DZ	0.25	0/635	0.50	0/848
52	BW	0.28	0/603	0.49	0/797
52	DW	0.28	0/603	0.49	0/797
All	All	0.27	13/306360 (0.0%)	0.70	91/457969 (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	15
1	CA	0	19
23	BB	0	37
23	DB	0	37
All	All	0	108

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	DB	1086	A	C5-C6	-16.20	1.26	1.41
23	BB	1086	A	C5-C6	-16.10	1.26	1.41
23	DB	1088	A	C6-N1	-10.51	1.28	1.35
23	BB	1088	A	C6-N1	-10.50	1.28	1.35
23	BB	1060	U	C2-N3	7.92	1.43	1.37
23	DB	1060	U	C2-N3	7.87	1.43	1.37
23	BB	1086	A	N3-C4	-7.12	1.30	1.34
23	DB	1086	A	N3-C4	-7.11	1.30	1.34
23	BB	1086	A	N7-C5	-6.38	1.35	1.39
23	DB	1086	A	N7-C5	-6.21	1.35	1.39
1	AA	1533	C	C4'-C3'	5.84	1.59	1.53
1	AA	495	A	N3-C4	-5.39	1.31	1.34
1	CA	495	A	N3-C4	-5.32	1.31	1.34

All (91) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DB	2204	G	O5'-P-OP1	-30.02	74.68	110.70
23	BB	2204	G	O5'-P-OP2	-28.49	76.52	110.70
23	DB	2791	G	O5'-P-OP2	-28.43	76.59	110.70
23	BB	2791	G	O5'-P-OP1	-27.42	77.80	110.70
23	DB	2204	G	O5'-P-OP2	17.65	131.88	110.70
23	DB	2791	G	O5'-P-OP1	17.55	131.76	110.70
23	BB	2204	G	O5'-P-OP1	17.49	131.69	110.70
23	BB	2791	G	O5'-P-OP2	17.14	131.27	110.70
23	DB	2790	U	OP2-P-O3'	15.09	138.40	105.20
23	DB	2203	U	OP1-P-O3'	14.30	136.65	105.20
23	BB	2203	U	OP2-P-O3'	13.96	135.91	105.20
23	BB	2790	U	OP1-P-O3'	13.80	135.56	105.20
23	BB	973	A	C5'-C4'-C3'	-8.67	102.13	116.00
23	DB	278	A	C5'-C4'-C3'	-8.66	102.15	116.00
23	BB	1552	A	N9-C1'-C2'	-8.41	102.75	112.00
23	DB	1552	A	N9-C1'-C2'	-8.40	102.76	112.00
23	BB	1088	A	N1-C6-N6	-8.22	113.67	118.60
23	DB	1088	A	N1-C6-N6	-8.16	113.70	118.60
23	BB	241	A	C5'-C4'-C3'	-8.07	103.09	116.00
1	AA	1499	A	C5'-C4'-O4'	-7.96	99.55	109.10
1	CA	765	G	N9-C1'-C2'	-7.36	103.90	112.00
1	AA	765	G	N9-C1'-C2'	-7.35	103.91	112.00
23	DB	1060	U	C5-C4-O4	-7.25	121.55	125.90
23	BB	1060	U	C5-C4-O4	-7.23	121.56	125.90
1	CA	576	C	O5'-P-OP1	-7.21	99.21	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DB	1086	A	C4-C5-C6	7.18	120.59	117.00
23	BB	1086	A	C4-C5-C6	7.16	120.58	117.00
1	CA	232	G	C5'-C4'-C3'	-6.90	104.95	116.00
1	AA	965	U	O4'-C4'-C3'	6.79	111.53	106.10
23	BB	2272	U	C5-C4-O4	-6.70	121.88	125.90
23	DB	2272	U	C5-C4-O4	-6.69	121.89	125.90
1	AA	232	G	C5'-C4'-C3'	-6.57	105.49	116.00
23	BB	1439	A	N9-C1'-C2'	-6.46	104.89	112.00
23	DB	1088	A	C5-C6-N6	6.43	128.85	123.70
1	AA	1126	U	C5'-C4'-C3'	-6.42	105.72	116.00
23	DB	1439	A	N9-C1'-C2'	-6.42	104.94	112.00
23	BB	268	C	C5'-C4'-C3'	-6.39	105.78	116.00
23	BB	1088	A	C5-C6-N6	6.38	128.81	123.70
22	BA	87	U	C4'-C3'-O3'	6.35	125.71	113.00
1	CA	95	C	C5'-C4'-C3'	-6.18	106.11	116.00
23	DB	2790	U	O3'-P-O5'	-6.17	92.27	104.00
23	DB	1086	A	C6-C5-N7	-6.17	127.98	132.30
23	BB	1086	A	C6-C5-N7	-6.15	128.00	132.30
1	AA	438	U	N1-C1'-C2'	-6.07	105.32	112.00
22	DA	87	U	C4'-C3'-O3'	6.04	125.08	113.00
1	CA	438	U	N1-C1'-C2'	-6.01	105.39	112.00
23	DB	2272	U	C5'-C4'-C3'	-5.99	106.42	116.00
1	AA	1361	G	N9-C1'-C2'	-5.96	105.44	112.00
23	DB	268	C	C5'-C4'-C3'	-5.90	106.56	116.00
23	DB	2108	A	N9-C1'-C2'	-5.86	105.56	112.00
23	DB	456	C	C5'-C4'-C3'	-5.85	106.64	116.00
1	CA	1300	G	C4'-C3'-C2'	5.84	108.44	102.60
23	BB	456	C	C5'-C4'-C3'	-5.77	106.76	116.00
23	BB	2191	A	C5'-C4'-C3'	-5.77	106.77	116.00
23	BB	1086	A	C2-N3-C4	-5.69	107.75	110.60
23	BB	1060	U	N1-C2-O2	-5.67	118.83	122.80
23	DB	1086	A	C2-N3-C4	-5.64	107.78	110.60
23	DB	1664	A	C5'-C4'-C3'	-5.63	107.00	116.00
23	BB	2894	G	N9-C1'-C2'	-5.62	105.81	112.00
23	DB	2894	G	N9-C1'-C2'	-5.62	105.82	112.00
23	DB	2203	U	O3'-P-O5'	-5.59	93.39	104.00
1	CA	1248	A	C5'-C4'-C3'	-5.58	107.06	116.00
1	CA	1249	C	C5'-C4'-C3'	-5.58	107.07	116.00
23	DB	1060	U	N1-C2-O2	-5.58	118.89	122.80
23	BB	2471	A	C5'-C4'-C3'	-5.50	107.20	116.00
23	BB	2272	U	C5'-C4'-C3'	-5.47	107.25	116.00
1	CA	1534	A	C2'-C3'-O3'	-5.42	97.57	109.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1213	A	C1'-O4'-C4'	-5.42	105.56	109.90
23	DB	1054	A	N9-C1'-C2'	-5.41	106.05	112.00
23	BB	1552	A	C4'-C3'-O3'	5.38	123.76	113.00
23	DB	1060	U	N3-C2-O2	5.36	125.95	122.20
23	BB	1060	U	N3-C2-O2	5.33	125.93	122.20
1	AA	86	G	N9-C1'-C2'	5.32	120.92	114.00
23	BB	560	C	C5'-C4'-C3'	-5.29	107.53	116.00
23	DB	1552	A	C4'-C3'-O3'	5.29	123.57	113.00
1	CA	1181	G	C5'-C4'-C3'	-5.26	107.58	116.00
1	AA	87	C	C4'-C3'-O3'	5.26	123.51	113.00
23	BB	2733	A	N9-C1'-C2'	-5.26	106.22	112.00
1	AA	1328	C	C5'-C4'-C3'	5.24	124.38	116.00
23	BB	2104	C	C4'-C3'-O3'	5.24	123.48	113.00
23	BB	2192	U	C5'-C4'-C3'	-5.22	107.65	116.00
23	DB	2471	A	C5'-C4'-C3'	-5.21	107.67	116.00
23	DB	2894	G	C5'-C4'-C3'	-5.19	107.70	116.00
1	CA	1023	U	C4'-C3'-O3'	5.18	123.36	113.00
23	BB	2203	U	O3'-P-O5'	-5.17	94.17	104.00
23	DB	2733	A	N9-C1'-C2'	-5.17	106.31	112.00
23	BB	143	C	C5'-C4'-C3'	-5.13	107.79	116.00
23	BB	479	A	C4'-C3'-O3'	-5.09	98.71	109.40
1	AA	1432	G	N9-C1'-C2'	-5.03	106.47	112.00
23	DB	2760	C	C5'-C4'-C3'	-5.03	107.96	116.00
23	DB	2619	C	C5'-C4'-C3'	-5.02	107.96	116.00

There are no chirality outliers.

All (108) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AA	1002	G	Sidechain
1	AA	1306	A	Sidechain
1	AA	1331	G	Sidechain
1	AA	1337	G	Sidechain
1	AA	1361	G	Sidechain
1	AA	1432	G	Sidechain
1	AA	1441	A	Sidechain
1	AA	187	G	Sidechain
1	AA	281	G	Sidechain
1	AA	437	U	Sidechain
1	AA	438	U	Sidechain
1	AA	450	G	Sidechain
1	AA	496	A	Sidechain

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Mol	Chain	Res	Type	Group
1	AA	703	G	Sidechain
1	AA	992	U	Sidechain
23	BB	1047	G	Sidechain
23	BB	1060	U	Sidechain
23	BB	1086	A	Sidechain
23	BB	1088	A	Sidechain
23	BB	1190	G	Sidechain
23	BB	1215	G	Sidechain
23	BB	1419	A	Sidechain
23	BB	1439	A	Sidechain
23	BB	1572	A	Sidechain
23	BB	1738	G	Sidechain
23	BB	1828	G	Sidechain
23	BB	2062	A	Sidechain
23	BB	2185	U	Sidechain
23	BB	221	A	Sidechain
23	BB	222	A	Sidechain
23	BB	2272	U	Sidechain
23	BB	232	G	Sidechain
23	BB	2454	G	Sidechain
23	BB	2471	A	Sidechain
23	BB	2503	A	Sidechain
23	BB	2508	G	Sidechain
23	BB	2638	G	Sidechain
23	BB	2733	A	Sidechain
23	BB	2770	G	Sidechain
23	BB	2848	G	Sidechain
23	BB	2868	A	Sidechain
23	BB	2883	A	Sidechain
23	BB	299	A	Sidechain
23	BB	3	U	Sidechain
23	BB	361	G	Sidechain
23	BB	370	G	Sidechain
23	BB	500	G	Sidechain
23	BB	51	G	Sidechain
23	BB	630	G	Sidechain
23	BB	633	A	Sidechain
23	BB	727	A	Sidechain
23	BB	757	G	Sidechain
1	CA	1012	A	Sidechain
1	CA	1027	C	Sidechain
1	CA	1033	G	Sidechain

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Mol	Chain	Res	Type	Group
1	CA	1133	G	Sidechain
1	CA	1144	G	Sidechain
1	CA	1313	U	Sidechain
1	CA	1331	G	Sidechain
1	CA	1356	G	Sidechain
1	CA	1432	G	Sidechain
1	CA	1441	A	Sidechain
1	CA	187	G	Sidechain
1	CA	281	G	Sidechain
1	CA	437	U	Sidechain
1	CA	438	U	Sidechain
1	CA	450	G	Sidechain
1	CA	496	A	Sidechain
1	CA	703	G	Sidechain
1	CA	82	G	Sidechain
1	CA	935	A	Sidechain
23	DB	1047	G	Sidechain
23	DB	1054	A	Sidechain
23	DB	1060	U	Sidechain
23	DB	1086	A	Sidechain
23	DB	1088	A	Sidechain
23	DB	1190	G	Sidechain
23	DB	1215	G	Sidechain
23	DB	1419	A	Sidechain
23	DB	1439	A	Sidechain
23	DB	1572	A	Sidechain
23	DB	1738	G	Sidechain
23	DB	1814	G	Sidechain
23	DB	1828	G	Sidechain
23	DB	2062	A	Sidechain
23	DB	2108	A	Sidechain
23	DB	221	A	Sidechain
23	DB	222	A	Sidechain
23	DB	2272	U	Sidechain
23	DB	232	G	Sidechain
23	DB	2454	G	Sidechain
23	DB	2471	A	Sidechain
23	DB	2503	A	Sidechain
23	DB	2508	G	Sidechain
23	DB	2638	G	Sidechain
23	DB	2733	A	Sidechain
23	DB	2770	G	Sidechain

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Mol	Chain	Res	Type	Group
23	DB	2848	G	Sidechain
23	DB	2868	A	Sidechain
23	DB	2883	A	Sidechain
23	DB	299	A	Sidechain
23	DB	370	G	Sidechain
23	DB	500	G	Sidechain
23	DB	51	G	Sidechain
23	DB	630	G	Sidechain
23	DB	633	A	Sidechain
23	DB	727	A	Sidechain
23	DB	757	G	Sidechain

5.2 Too-close contacts [i](#)

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	32831	0	16521	1458	0
1	CA	32831	0	16521	1414	0
2	AC	1624	0	1699	205	0
2	CC	1624	0	1699	191	0
3	AD	1643	0	1710	179	0
3	CD	1643	0	1710	177	0
4	AE	1105	0	1148	129	0
4	CE	1105	0	1148	121	0
5	AF	817	0	808	89	0
5	CF	817	0	808	91	0
6	AG	1174	0	1230	146	0
6	CG	1196	0	1246	133	0
7	AH	979	0	1034	89	0
7	CH	979	0	1034	91	0
8	AI	1022	0	1070	180	0
8	CI	1022	0	1070	146	0
9	AJ	786	0	828	85	0
9	CJ	786	0	828	103	0
10	AK	877	0	887	110	0
10	CK	877	0	887	100	0
11	AL	955	0	1019	96	0
11	CL	955	0	1019	97	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	AM	883	0	944	135	0
12	CM	876	0	937	138	0
13	AP	649	0	666	65	0
13	CP	638	0	656	66	0
14	AQ	648	0	691	63	0
14	CQ	657	0	702	62	0
15	AR	455	0	478	35	0
15	CR	455	0	478	37	0
16	AS	637	0	665	97	0
16	CS	644	0	675	115	0
17	AT	665	0	714	60	0
17	CT	665	0	714	61	0
18	AB	1704	0	1732	209	0
18	CB	1704	0	1732	229	0
19	AU	425	0	449	57	0
19	CU	425	0	449	54	0
20	AO	714	0	734	63	0
20	CO	714	0	734	62	0
21	AN	774	0	827	102	0
21	CN	774	0	827	114	0
22	BA	2507	0	1270	116	0
22	DA	2507	0	1270	111	0
23	BB	60995	0	30679	2412	0
23	DB	60995	0	30678	2455	0
24	BI	1032	0	1088	109	0
24	DI	1032	0	1088	168	0
25	BC	2082	0	2157	234	0
25	DC	2082	0	2157	239	0
26	BD	1565	0	1616	234	0
26	DD	1565	0	1616	239	0
27	BK	930	0	1000	153	0
27	DK	930	0	1000	154	0
28	BP	917	0	965	126	0
28	DP	917	0	965	132	0
29	BE	1552	0	1619	208	0
29	DE	1552	0	1619	202	0
30	BY	449	0	491	59	0
30	DY	449	0	491	53	0
31	B0	444	0	461	48	0
31	D0	444	0	461	44	0
32	B4	302	0	340	42	0
32	D4	302	0	340	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
33	B1	409	0	440	54	0
33	D1	409	0	440	50	0
34	B3	504	0	574	47	0
34	D3	504	0	574	52	0
35	BV	753	0	780	83	0
35	DV	753	0	780	86	0
36	B2	377	0	418	37	0
36	D2	377	0	418	43	0
37	BL	1045	0	1117	138	0
37	DL	1045	0	1117	144	0
38	BM	1074	0	1157	123	0
38	DM	1074	0	1157	119	0
39	BX	509	0	543	46	0
39	DX	509	0	543	50	0
40	BH	1111	0	1148	172	0
40	DH	1111	0	1148	147	0
41	BJ	1129	0	1162	134	0
41	DJ	1129	0	1162	141	0
42	BN	960	0	1000	137	0
42	DN	960	0	1000	133	0
43	BO	892	0	923	91	0
43	DO	892	0	923	94	0
44	BQ	947	0	1022	150	0
44	DQ	947	0	1022	143	0
45	BS	857	0	922	97	0
45	DS	857	0	922	98	0
46	BU	779	0	834	116	0
46	DU	779	0	834	114	0
47	BF	1420	0	1460	264	0
47	DF	1420	0	1460	249	0
48	BG	1323	0	1374	187	0
48	DG	1323	0	1374	178	0
49	BR	816	0	839	105	0
49	DR	816	0	839	112	0
50	BT	738	0	807	115	0
50	DT	738	0	807	110	0
51	BZ	625	0	652	75	0
51	DZ	625	0	652	71	0
52	BW	596	0	610	122	0
52	DW	596	0	610	130	0
53	AA	60	0	0	0	0
53	BB	110	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
53	CA	58	0	0	0	0
53	CE	1	0	0	0	0
53	DB	110	0	0	0	0
53	DN	1	0	0	0	0
54	AA	23	0	24	2	0
54	CA	23	0	24	1	0
55	B4	1	0	0	0	0
55	D4	1	0	0	0	0
56	AA	288	0	0	6	0
56	AE	3	0	0	1	0
56	AK	1	0	0	0	0
56	AL	4	0	0	0	0
56	AN	2	0	0	0	0
56	AP	1	0	0	0	0
56	AT	1	0	0	0	0
56	BB	494	0	0	4	0
56	BC	4	0	0	0	0
56	BE	3	0	0	0	0
56	BH	1	0	0	0	0
56	BL	4	0	0	0	0
56	BT	1	0	0	0	0
56	CA	275	0	0	4	0
56	CE	4	0	0	0	0
56	CK	1	0	0	0	0
56	CL	5	0	0	0	0
56	CN	5	0	0	0	0
56	CP	1	0	0	0	0
56	CT	2	0	0	0	0
56	DB	500	0	0	9	0
56	DC	3	0	0	0	0
56	DD	1	0	0	0	0
56	DE	1	0	0	0	0
56	DJ	1	0	0	0	0
56	DL	3	0	0	0	0
56	DN	2	0	0	0	0
56	DP	1	0	0	0	0
56	DR	1	0	0	0	0
All	All	284033	0	190711	17874	0

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 38.

All (17874) 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)
23:DB:1099:G:H8	24:DI:3:LYS:N	1.36	1.21
23:BB:855:G:H21	52:BW:23:LYS:HG2	1.11	1.13
23:DB:322:A:H5'	23:DB:340:A:H1'	1.32	1.12
2:AC:70:ALA:HA	2:AC:105:VAL:HG21	1.26	1.11
23:BB:1205:A:H62	29:BE:165:HIS:HB2	1.11	1.10
23:BB:1244:A:H5''	37:BL:8:PRO:HD3	1.35	1.08
23:DB:1244:A:H5''	37:DL:8:PRO:HD3	1.32	1.08
2:CC:52:SER:HB3	2:CC:114:LEU:HG	1.37	1.07
23:BB:322:A:H5'	23:BB:340:A:H1'	1.28	1.07
23:DB:1099:G:C8	24:DI:3:LYS:N	2.25	1.05
13:AP:28:ARG:HD2	13:AP:29:ASN:H	1.21	1.04
40:DH:31:VAL:HB	40:DH:32:PRO:HD3	1.39	1.04
9:AJ:7:ARG:HB2	9:AJ:101:SER:H	1.23	1.03
38:BM:59:ARG:HE	38:BM:60:GLN:HG2	1.22	1.03
25:DC:144:GLU:HG3	25:DC:151:GLY:H	1.23	1.03
40:BH:84:ALA:HA	40:BH:90:LEU:HA	1.39	1.03
38:DM:41:LEU:HB2	38:DM:94:ALA:HB3	1.40	1.03
29:BE:61:ARG:HH12	29:BE:64:GLY:HA3	1.23	1.03
23:DB:1098:A:H3'	24:DI:3:LYS:HA	1.41	1.03
23:DB:1099:G:O5'	24:DI:4:VAL:N	1.92	1.03
40:BH:31:VAL:HB	40:BH:32:PRO:HD3	1.39	1.02
18:CB:46:VAL:HG13	18:CB:47:PRO:HD3	1.42	1.02
1:AA:71:A:H61	1:AA:99:C:H1'	1.23	1.02
13:CP:28:ARG:HD2	13:CP:29:ASN:H	1.21	1.01
1:AA:1313:U:H5''	16:AS:5:LYS:HB3	1.39	1.01
51:BZ:45:ARG:HE	51:BZ:47:VAL:HG12	1.18	1.01
52:BW:17:ALA:HA	52:BW:35:ILE:HG23	1.39	1.00
38:DM:59:ARG:HE	38:DM:60:GLN:HG2	1.23	1.00
1:AA:1211:U:H4'	1:AA:1213:A:H1'	1.44	1.00
23:BB:674:G:H5''	29:BE:71:GLY:H	1.20	1.00
18:CB:40:ILE:HD13	18:CB:201:GLY:HA2	1.44	0.99
12:AM:47:LEU:HG	12:AM:52:ILE:HD13	1.44	0.99
50:DT:55:VAL:HA	50:DT:87:LEU:HA	1.44	0.99
9:AJ:36:VAL:HA	9:AJ:76:ILE:HG22	1.38	0.99
23:BB:674:G:H1'	29:BE:69:ARG:HE	1.26	0.99
50:BT:55:VAL:HA	50:BT:87:LEU:HA	1.44	0.99
38:BM:41:LEU:HB2	38:BM:94:ALA:HB3	1.43	0.99
21:CN:51:PRO:HB2	21:CN:54:SER:HB3	1.45	0.99
45:DS:70:LYS:HD3	45:DS:110:ARG:HA	1.45	0.99
1:AA:1048:G:H21	1:AA:1214:C:H5	1.10	0.98
6:AG:107:ALA:HB1	6:AG:119:LEU:HA	1.45	0.98
25:DC:244:VAL:HB	25:DC:249:VAL:H	1.26	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:10:GLY:HA3	26:DD:26:VAL:H	1.27	0.98
40:DH:131:SER:HB2	40:DH:141:LYS:HA	1.45	0.98
21:AN:50:LEU:H	21:AN:51:PRO:HD2	1.28	0.98
1:CA:1493:A:HO2'	1:CA:1494:G:H8	1.02	0.98
52:DW:17:ALA:HA	52:DW:35:ILE:HG23	1.45	0.98
40:BH:47:PHE:HA	40:BH:50:ARG:HE	1.26	0.97
18:CB:53:LEU:HD13	18:CB:216:VAL:HG12	1.46	0.97
23:DB:1098:A:H3'	24:DI:3:LYS:CA	1.93	0.97
18:AB:202:ASN:HD22	18:AB:204:ASP:H	0.98	0.97
44:DQ:63:ARG:HH12	44:DQ:96:ASP:HB2	1.30	0.97
25:DC:144:GLU:HA	25:DC:151:GLY:HA2	1.45	0.97
25:BC:144:GLU:HG3	25:BC:151:GLY:H	1.26	0.97
26:BD:10:GLY:HA3	26:BD:26:VAL:H	1.29	0.96
2:CC:59:PRO:HG2	2:CC:62:SER:HB2	1.47	0.96
51:DZ:45:ARG:HE	51:DZ:47:VAL:HG12	1.26	0.96
45:BS:70:LYS:HD3	45:BS:110:ARG:HA	1.47	0.96
16:CS:39:ILE:HD13	16:CS:65:MET:HG2	1.46	0.96
15:AR:72:ARG:H	15:AR:72:ARG:HH11	1.04	0.96
52:BW:43:LYS:HD2	52:BW:79:ILE:HD11	1.46	0.96
23:DB:275:C:H2'	23:DB:276:U:H4'	1.45	0.96
27:DK:35:VAL:HG23	27:DK:36:GLY:H	1.29	0.96
15:CR:72:ARG:H	15:CR:72:ARG:HH11	1.02	0.95
51:BZ:71:LEU:HD13	51:BZ:76:GLU:HB3	1.48	0.95
40:BH:80:ILE:H	40:BH:144:VAL:HG22	1.32	0.95
2:CC:190:THR:HG23	2:CC:192:TYR:H	1.28	0.95
29:DE:61:ARG:HH12	29:DE:64:GLY:HA3	1.29	0.95
25:BC:144:GLU:HA	25:BC:151:GLY:HA2	1.47	0.95
12:CM:21:ILE:HB	12:CM:24:VAL:HG22	1.48	0.95
51:DZ:71:LEU:HD13	51:DZ:76:GLU:HB3	1.49	0.95
9:AJ:53:ILE:HG22	9:AJ:61:ALA:HB1	1.47	0.95
8:CI:51:LEU:HA	8:CI:54:VAL:HG22	1.49	0.95
23:DB:1099:G:H8	24:DI:3:LYS:H	0.99	0.95
25:BC:244:VAL:HB	25:BC:249:VAL:H	1.27	0.94
27:BK:35:VAL:HG23	27:BK:36:GLY:H	1.28	0.94
6:AG:14:ASP:HB3	6:AG:19:SER:H	1.29	0.94
18:AB:209:VAL:HG23	18:AB:210:THR:H	1.30	0.94
52:DW:37:VAL:HG12	52:DW:38:ARG:H	1.32	0.94
40:DH:125:THR:HA	40:DH:146:VAL:HB	1.46	0.94
19:AU:36:PHE:HB3	19:AU:40:PRO:HD3	1.50	0.94
6:CG:75:LYS:HZ2	6:CG:76:SER:H	1.15	0.94
8:AI:29:ILE:HA	8:AI:64:ILE:HB	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:120:A:H2'	1:AA:121:U:H5''	1.50	0.94
1:CA:1221:G:H4'	16:CS:76:THR:HG21	1.50	0.94
6:CG:78:ARG:HG2	6:CG:83:THR:HG22	1.49	0.94
42:DN:37:THR:HG22	42:DN:39:PRO:HD2	1.48	0.94
42:BN:37:THR:HG22	42:BN:39:PRO:HD2	1.49	0.93
18:CB:56:LEU:HD22	18:CB:216:VAL:HB	1.48	0.93
42:DN:101:GLY:HA2	42:DN:110:MET:H	1.32	0.93
23:BB:2134:A:H62	23:BB:2157:G:H21	1.14	0.93
1:AA:1361:G:H2'	1:AA:1362:A:H5''	1.48	0.93
23:BB:1597:A:H5''	23:BB:1598:A:H5'	1.51	0.93
44:BQ:63:ARG:HH12	44:BQ:96:ASP:HB2	1.34	0.93
3:AD:22:SER:H	3:AD:109:THR:HG22	1.30	0.93
2:CC:155:ARG:H	2:CC:162:ALA:HA	1.31	0.93
23:DB:877:A:H61	23:DB:898:C:H2'	1.32	0.93
37:DL:23:ILE:HD12	37:DL:23:ILE:H	1.33	0.93
8:CI:117:LEU:HD22	8:CI:123:ARG:HB3	1.50	0.93
16:CS:18:VAL:HG21	16:CS:43:MET:HG2	1.49	0.93
23:BB:876:C:H41	23:BB:901:C:N4	1.67	0.93
23:DB:1597:A:H5''	23:DB:1598:A:H5'	1.50	0.92
2:AC:78:LYS:HG2	2:AC:81:GLU:HB2	1.51	0.92
3:CD:22:SER:H	3:CD:109:THR:HG22	1.31	0.92
19:CU:36:PHE:HB3	19:CU:40:PRO:HD3	1.51	0.92
23:BB:858:G:N3	23:BB:2268:A:H2'	1.82	0.92
23:DB:45:G:H5'	23:DB:46:G:H5'	1.51	0.92
42:BN:101:GLY:HA2	42:BN:110:MET:H	1.33	0.92
1:CA:120:A:H2'	1:CA:121:U:H5''	1.48	0.92
11:AL:23:LEU:HG	11:AL:24:GLU:HG3	1.49	0.92
13:AP:66:THR:HG22	13:AP:67:ILE:H	1.35	0.92
2:CC:26:LYS:HG3	2:CC:27:GLU:H	1.33	0.92
23:DB:704:G:H2'	23:DB:726:G:H22	1.34	0.92
26:BD:148:GLN:HB2	26:BD:152:PRO:HG2	1.52	0.92
1:CA:812:G:HO2'	1:CA:813:U:H6	0.98	0.92
43:DO:5:SER:HA	43:DO:8:ILE:HD12	1.52	0.92
25:BC:146:LYS:HB3	25:BC:147:PRO:HD2	1.51	0.92
37:DL:123:ARG:HA	37:DL:143:GLU:HB3	1.50	0.92
3:CD:152:SER:HA	3:CD:155:LYS:HD3	1.53	0.91
12:CM:79:LEU:HD22	12:CM:86:ARG:HE	1.34	0.91
43:DO:66:GLY:HA3	43:DO:102:ARG:HH21	1.34	0.91
9:AJ:8:ILE:HA	9:AJ:100:ILE:HG22	1.49	0.91
23:BB:45:G:H5'	23:BB:46:G:H5'	1.50	0.91
37:BL:123:ARG:HA	37:BL:143:GLU:HB3	1.50	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:69:ALA:HA	40:BH:140:ALA:HA	1.50	0.91
2:AC:26:LYS:HG3	2:AC:27:GLU:H	1.36	0.91
11:CL:23:LEU:HG	11:CL:24:GLU:HG3	1.53	0.90
23:DB:27:G:H22	23:DB:512:G:H2'	1.35	0.90
4:CE:81:GLN:H	4:CE:146:MET:HE3	1.36	0.90
43:BO:66:GLY:HA3	43:BO:102:ARG:HH21	1.37	0.90
23:DB:1205:A:H62	29:DE:165:HIS:HB2	1.31	0.90
23:DB:858:G:N3	23:DB:2268:A:H2'	1.86	0.90
3:AD:152:SER:HA	3:AD:155:LYS:HD3	1.53	0.90
8:CI:5:TYR:HB2	8:CI:20:ILE:HB	1.51	0.90
16:CS:39:ILE:HB	16:CS:66:VAL:HA	1.52	0.90
33:D1:33:LEU:HB3	33:D1:51:ALA:HB3	1.51	0.90
23:DB:276:U:H2'	23:DB:278:A:H62	1.36	0.90
23:DB:1652:A:H62	42:DN:11:ASN:HD21	1.19	0.90
32:B4:7:VAL:HG13	32:B4:8:LYS:H	1.35	0.90
23:BB:674:G:H5''	29:BE:71:GLY:N	1.85	0.90
18:CB:159:ALA:HB1	18:CB:183:PHE:HE1	1.33	0.90
5:AF:92:THR:HG22	5:AF:94:HIS:H	1.36	0.90
25:BC:117:SER:HB2	25:BC:128:THR:HB	1.52	0.90
43:BO:5:SER:HA	43:BO:8:ILE:HD12	1.54	0.90
6:CG:14:ASP:H	6:CG:19:SER:H	1.17	0.90
52:BW:37:VAL:HG12	52:BW:38:ARG:H	1.35	0.90
18:CB:156:LEU:HD23	18:CB:180:ILE:HD11	1.54	0.90
13:CP:4:ILE:HB	13:CP:67:ILE:HD12	1.53	0.90
25:DC:93:VAL:HG13	25:DC:94:LEU:H	1.35	0.89
48:DG:85:LYS:HA	48:DG:131:VAL:HG12	1.54	0.89
32:D4:7:VAL:HG13	32:D4:8:LYS:H	1.36	0.89
1:AA:79:G:H2'	1:AA:80:A:H8	1.36	0.89
23:BB:1060:U:N3	23:BB:1088:A:N7	2.20	0.89
48:BG:85:LYS:HA	48:BG:131:VAL:HG12	1.54	0.89
1:AA:1086:U:H3	1:AA:1099:G:H22	1.17	0.89
50:DT:15:HIS:H	50:DT:32:LEU:HA	1.37	0.89
50:BT:57:VAL:HG22	50:BT:58:VAL:H	1.35	0.89
13:CP:66:THR:HG22	13:CP:67:ILE:H	1.37	0.89
23:DB:1060:U:N3	23:DB:1088:A:N7	2.21	0.89
37:DL:29:LYS:HG3	37:DL:30:THR:HG23	1.55	0.89
52:DW:43:LYS:HD2	52:DW:79:ILE:HD11	1.52	0.89
4:CE:93:VAL:HG13	4:CE:126:ALA:HB2	1.53	0.89
33:D1:49:LYS:HG2	33:D1:50:GLU:H	1.37	0.89
25:DC:41:GLY:HA3	25:DC:53:ILE:HG21	1.55	0.89
26:DD:148:GLN:HB2	26:DD:152:PRO:HG2	1.53	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:36:VAL:HG13	9:CJ:76:ILE:HA	1.54	0.89
41:DJ:24:THR:HA	41:DJ:63:ALA:HB3	1.55	0.89
1:AA:1057:G:H4'	2:AC:196:GLY:H	1.36	0.89
42:DN:41:ALA:HB1	42:DN:113:ILE:HD13	1.52	0.89
52:DW:51:GLY:HA3	52:DW:59:PHE:HB2	1.55	0.89
23:BB:855:G:N2	52:BW:23:LYS:HG2	1.87	0.88
52:BW:39:GLN:HG3	52:BW:42:THR:HB	1.55	0.88
25:BC:93:VAL:HG13	25:BC:94:LEU:H	1.37	0.88
4:CE:113:VAL:HG23	4:CE:114:LEU:H	1.37	0.88
5:CF:92:THR:HG22	5:CF:94:HIS:H	1.36	0.88
33:D1:26:LYS:HD3	33:D1:52:LYS:HB3	1.55	0.88
23:DB:845:A:H2'	23:DB:846:U:H5''	1.55	0.88
25:DC:117:SER:HB2	25:DC:128:THR:HB	1.55	0.88
29:DE:58:LYS:HE2	29:DE:60:TRP:HD1	1.35	0.88
1:AA:1057:G:H5''	2:AC:153:SER:HB2	1.53	0.88
40:BH:99:ILE:HG21	40:BH:130:VAL:HB	1.56	0.88
21:CN:50:LEU:H	21:CN:51:PRO:HD2	1.39	0.88
44:DQ:105:PHE:HA	44:DQ:108:LEU:HD12	1.54	0.88
24:BI:129:GLU:HB3	24:BI:133:ARG:HH12	1.36	0.88
37:BL:116:VAL:HG13	37:BL:117:THR:H	1.38	0.88
1:CA:981:U:H4'	21:CN:60:ARG:HD2	1.54	0.88
23:DB:547:A:H2'	23:DB:547:A:N3	1.89	0.88
23:BB:704:G:H2'	23:BB:726:G:H22	1.38	0.88
24:DI:11:GLN:HG2	24:DI:55:PRO:HB3	1.56	0.88
1:AA:1328:C:H5''	12:AM:27:THR:HG21	1.55	0.88
21:AN:60:ARG:HG3	21:AN:62:ARG:HG2	1.55	0.88
42:DN:2:ARG:HG2	42:DN:5:LYS:HB2	1.55	0.88
1:AA:1060:U:H4'	9:AJ:54:SER:HB2	1.56	0.88
13:AP:4:ILE:HB	13:AP:67:ILE:HD12	1.54	0.88
37:BL:23:ILE:HD12	37:BL:23:ILE:H	1.38	0.88
13:CP:4:ILE:HG12	13:CP:21:VAL:HG22	1.54	0.88
25:DC:146:LYS:HB3	25:DC:147:PRO:HD2	1.52	0.88
12:AM:3:ILE:HA	12:AM:56:ARG:HD3	1.53	0.88
33:B1:26:LYS:HD3	33:B1:52:LYS:HB3	1.54	0.88
28:DP:56:SER:HB2	28:DP:75:THR:HB	1.56	0.88
1:AA:664:G:H22	1:AA:741:G:H1	1.22	0.88
41:BJ:24:THR:HA	41:BJ:63:ALA:HB3	1.54	0.88
23:BB:27:G:H22	23:BB:512:G:H2'	1.38	0.87
25:BC:41:GLY:HA3	25:BC:53:ILE:HG21	1.54	0.87
44:BQ:105:PHE:HA	44:BQ:108:LEU:HD12	1.55	0.87
50:BT:15:HIS:H	50:BT:32:LEU:HA	1.39	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:64:VAL:O	25:DC:65:ASP:HB3	1.74	0.87
1:AA:79:G:H2'	1:AA:80:A:C8	2.08	0.87
6:AG:112:ASP:HB2	6:AG:118:ARG:HG2	1.55	0.87
38:DM:19:GLY:HA2	38:DM:98:PRO:HD2	1.54	0.87
33:B1:49:LYS:HG2	33:B1:50:GLU:H	1.38	0.87
43:BO:64:TYR:HB3	43:BO:67:ASN:HB2	1.55	0.87
6:CG:145:GLU:HA	6:CG:148:LYS:HB2	1.55	0.87
12:CM:28:ARG:HH21	12:CM:62:PHE:HB2	1.38	0.87
4:AE:14:LEU:HD23	4:AE:36:THR:HG22	1.55	0.87
6:AG:58:LEU:HD23	6:AG:58:LEU:H	1.40	0.87
19:AU:34:ARG:HD3	19:AU:39:LYS:HE3	1.57	0.87
23:DB:855:G:H21	52:DW:23:LYS:HG2	1.38	0.87
23:BB:1251:C:H2'	44:BQ:5:ARG:HH12	1.39	0.87
1:CA:1060:U:H4'	9:CJ:54:SER:HB2	1.55	0.87
37:DL:78:ARG:HB3	37:DL:113:ALA:HB2	1.57	0.87
50:DT:57:VAL:HG22	50:DT:58:VAL:H	1.36	0.87
4:AE:81:GLN:H	4:AE:146:MET:HE3	1.38	0.87
23:BB:845:A:H2'	23:BB:846:U:H5''	1.57	0.87
22:DA:2:G:H2'	22:DA:3:C:C6	2.09	0.87
52:DW:39:GLN:HG3	52:DW:42:THR:HB	1.56	0.87
18:AB:46:VAL:HG13	18:AB:47:PRO:HD3	1.57	0.87
2:AC:149:LYS:HB3	2:AC:200:TRP:HB2	1.57	0.87
21:AN:17:ASP:HA	21:AN:21:ALA:HB2	1.56	0.87
33:B1:33:LEU:HB3	33:B1:51:ALA:HB3	1.53	0.86
25:BC:64:VAL:O	25:BC:65:ASP:HB3	1.73	0.86
29:BE:52:VAL:HG12	29:BE:53:THR:H	1.40	0.86
2:AC:123:LEU:HA	2:AC:127:VAL:HG22	1.57	0.86
4:AE:113:VAL:HG23	4:AE:114:LEU:H	1.39	0.86
30:BY:12:ALA:HA	30:BY:15:ARG:HD3	1.57	0.86
23:DB:142:A:H2'	23:DB:143:C:C6	2.10	0.86
40:BH:116:ARG:HH12	40:BH:133:GLN:HB2	1.41	0.86
42:BN:41:ALA:HB1	42:BN:113:ILE:HD13	1.58	0.86
16:CS:39:ILE:HD11	16:CS:68:HIS:HB2	1.55	0.86
29:DE:52:VAL:HG12	29:DE:53:THR:H	1.39	0.86
50:DT:5:GLU:HA	50:DT:8:LEU:HB2	1.57	0.86
5:AF:38:ARG:HD3	5:AF:97:THR:HA	1.55	0.86
24:BI:27:LEU:HD23	24:BI:27:LEU:H	1.40	0.86
49:DR:39:LEU:HA	49:DR:53:PHE:HA	1.57	0.86
8:AI:20:ILE:HD13	8:AI:85:ALA:HB3	1.57	0.86
47:BF:134:GLN:H	47:BF:150:GLY:H	1.23	0.86
38:BM:19:GLY:HA2	38:BM:98:PRO:HD2	1.56	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:11:LEU:HD22	50:BT:11:LEU:H	1.39	0.86
1:CA:279:A:H5''	1:CA:280:C:H3'	1.58	0.86
15:CR:72:ARG:H	15:CR:72:ARG:NH1	1.73	0.86
4:AE:93:VAL:HG13	4:AE:126:ALA:HB2	1.55	0.86
6:CG:26:VAL:HG12	6:CG:42:VAL:HG11	1.58	0.86
6:CG:87:PRO:HG3	6:CG:148:LYS:HA	1.55	0.86
27:DK:47:ILE:HG12	27:DK:48:PRO:HD2	1.56	0.86
1:AA:1187:G:H21	21:AN:99:SER:HB2	1.40	0.86
1:CA:560:A:H4'	1:CA:561:U:H5''	1.58	0.86
22:DA:98:G:H1	35:DV:14:LYS:HB2	1.40	0.86
47:DF:134:GLN:H	47:DF:150:GLY:H	1.24	0.86
28:DP:21:PRO:HG2	28:DP:91:VAL:HG21	1.58	0.86
8:AI:12:LYS:H	8:AI:105:ARG:HH21	1.23	0.86
29:BE:111:GLU:HA	29:BE:114:ARG:HG3	1.58	0.86
18:CB:163:ILE:HG23	18:CB:164:ASP:H	1.39	0.86
37:DL:116:VAL:HG13	37:DL:117:THR:H	1.40	0.86
42:DN:2:ARG:HA	42:DN:5:LYS:HD3	1.57	0.86
29:BE:58:LYS:NZ	29:BE:58:LYS:H	1.74	0.85
47:BF:84:ILE:HG23	47:BF:85:GLY:H	1.41	0.85
1:CA:664:G:H22	1:CA:741:G:H1	1.24	0.85
9:CJ:35:GLN:HG3	9:CJ:77:VAL:HB	1.55	0.85
44:DQ:7:VAL:HA	44:DQ:10:ARG:HG2	1.56	0.85
29:BE:5:LEU:HD12	29:BE:10:SER:HB2	1.56	0.85
29:BE:58:LYS:HE2	29:BE:60:TRP:HD1	1.41	0.85
27:BK:47:ILE:HG12	27:BK:48:PRO:HD2	1.57	0.85
16:CS:51:HIS:HA	16:CS:56:HIS:HA	1.58	0.85
38:DM:35:ALA:HB2	38:DM:100:LYS:HB2	1.57	0.85
2:AC:76:ILE:HA	2:AC:83:VAL:HG23	1.56	0.85
1:CA:1171:A:H2'	1:CA:1172:C:C6	2.11	0.85
5:CF:38:ARG:HD3	5:CF:97:THR:HA	1.56	0.85
48:DG:30:GLY:HA3	48:DG:78:VAL:HA	1.56	0.85
37:DL:143:GLU:HG2	37:DL:144:GLU:H	1.38	0.85
9:AJ:57:VAL:HG22	9:AJ:58:ASN:H	1.39	0.85
23:BB:923:G:H1'	52:BW:23:LYS:HZ2	1.42	0.85
45:DS:26:GLY:H	45:DS:71:VAL:HG13	1.41	0.85
14:AQ:25:GLU:HA	14:AQ:40:THR:HA	1.57	0.85
2:CC:18:ASN:HA	2:CC:55:VAL:HG22	1.56	0.85
44:DQ:14:LYS:HA	44:DQ:17:LEU:HB3	1.57	0.85
46:DU:3:LYS:HB3	46:DU:82:VAL:HG21	1.57	0.85
49:BR:59:ILE:HG12	49:BR:101:ILE:HD13	1.57	0.85
23:BB:597:G:H21	37:BL:12:SER:HA	1.41	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:52:LEU:HG	9:CJ:62:ARG:HE	1.42	0.85
11:CL:24:GLU:HB2	11:CL:26:CYS:SG	2.17	0.85
27:DK:25:LEU:HD21	27:DK:40:LYS:HB2	1.58	0.85
3:AD:173:ASP:HB3	3:AD:178:GLU:HB2	1.57	0.85
1:AA:939:G:H5'	6:AG:101:ARG:HH12	1.41	0.85
9:CJ:6:ILE:HB	9:CJ:76:ILE:HG13	1.56	0.85
51:DZ:4:VAL:HG12	51:DZ:11:ARG:HG2	1.59	0.85
50:BT:29:THR:HA	50:BT:86:THR:HA	1.57	0.85
3:CD:173:ASP:HB3	3:CD:178:GLU:HB2	1.58	0.85
4:CE:14:LEU:HD23	4:CE:36:THR:HG22	1.58	0.85
29:DE:126:VAL:HG22	29:DE:127:GLU:H	1.41	0.85
37:BL:143:GLU:HG2	37:BL:144:GLU:H	1.40	0.84
7:CH:86:LYS:HB3	7:CH:90:GLU:HB3	1.59	0.84
23:DB:1579:A:H2'	23:DB:1580:A:C8	2.12	0.84
23:DB:2799:A:H4'	23:DB:2800:A:C8	2.12	0.84
50:DT:11:LEU:H	50:DT:11:LEU:HD22	1.39	0.84
21:AN:50:LEU:HD23	21:AN:51:PRO:HD3	1.58	0.84
42:BN:19:ALA:HA	42:BN:22:ARG:HB3	1.57	0.84
46:BU:3:LYS:HB3	46:BU:82:VAL:HG21	1.58	0.84
1:CA:1493:A:H3'	1:CA:1493:A:N3	1.92	0.84
8:CI:49:GLN:HA	8:CI:52:GLU:HG3	1.59	0.84
51:DZ:7:VAL:HG13	51:DZ:8:THR:HG23	1.59	0.84
4:AE:71:ILE:HG12	4:AE:72:ASN:H	1.40	0.84
12:CM:3:ILE:HG12	12:CM:52:ILE:HD11	1.60	0.84
22:BA:32:U:H4'	22:BA:52:A:H62	1.41	0.84
27:BK:25:LEU:HD21	27:BK:40:LYS:HB2	1.57	0.84
37:BL:78:ARG:HB3	37:BL:113:ALA:HB2	1.57	0.84
11:CL:78:VAL:HG12	11:CL:101:LEU:HD13	1.58	0.84
23:DB:1099:G:H5''	24:DI:3:LYS:N	1.92	0.84
42:DN:19:ALA:HA	42:DN:22:ARG:HB3	1.59	0.84
1:AA:922:G:H2'	1:AA:923:A:C8	2.11	0.84
4:CE:71:ILE:HG12	4:CE:72:ASN:H	1.40	0.84
44:BQ:14:LYS:HA	44:BQ:17:LEU:HB3	1.58	0.84
1:CA:17:U:H2'	1:CA:18:C:C6	2.13	0.84
4:CE:132:PRO:HG2	4:CE:133:ILE:HD12	1.60	0.84
41:BJ:17:VAL:HG23	41:BJ:137:PRO:HB2	1.58	0.84
1:AA:560:A:H4'	1:AA:561:U:H5''	1.59	0.84
10:AK:105:ARG:HH21	19:AU:10:PRO:HB3	1.43	0.84
52:BW:51:GLY:HA3	52:BW:59:PHE:HB2	1.57	0.84
12:CM:28:ARG:HH12	12:CM:32:ILE:HD12	1.41	0.84
1:AA:1125:U:H2'	1:AA:1126:U:H5''	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:202:ASN:ND2	18:AB:204:ASP:H	1.76	0.84
1:CA:920:U:H2'	1:CA:921:U:C6	2.13	0.84
15:AR:72:ARG:H	15:AR:72:ARG:NH1	1.75	0.84
23:BB:1310:G:H21	23:BB:1610:A:H8	1.26	0.84
1:CA:942:G:H21	8:CI:125:GLN:HE22	1.23	0.84
19:CU:34:ARG:HD3	19:CU:39:LYS:HE3	1.60	0.84
8:AI:20:ILE:HA	8:AI:62:LEU:HD12	1.59	0.83
48:BG:30:GLY:HA3	48:BG:78:VAL:HA	1.59	0.83
5:CF:3:HIS:HB2	5:CF:92:THR:HA	1.60	0.83
43:DO:64:TYR:HB3	43:DO:67:ASN:HB2	1.57	0.83
12:AM:14:ALA:HB2	12:AM:42:VAL:HG23	1.59	0.83
52:BW:39:GLN:HG2	52:BW:40:ARG:N	1.91	0.83
6:AG:71:THR:HA	6:AG:90:VAL:HG22	1.60	0.83
44:BQ:7:VAL:HA	44:BQ:10:ARG:HG2	1.58	0.83
23:DB:1310:G:H21	23:DB:1610:A:H8	1.25	0.83
30:DY:12:ALA:HA	30:DY:15:ARG:HD3	1.59	0.83
1:AA:973:G:H3'	1:AA:974:A:H5''	1.59	0.83
23:BB:364:C:H2'	23:BB:365:U:C6	2.14	0.83
27:BK:107:LEU:HD12	27:BK:107:LEU:H	1.43	0.83
14:CQ:25:GLU:HA	14:CQ:40:THR:HA	1.61	0.83
25:DC:94:LEU:HD13	25:DC:100:ARG:HD2	1.58	0.83
27:DK:105:ARG:H	27:DK:105:ARG:HD3	1.43	0.83
27:DK:54:LYS:H	27:DK:54:LYS:HD2	1.42	0.83
8:AI:51:LEU:HB2	8:AI:56:MET:SD	2.17	0.83
10:CK:33:ILE:HB	10:CK:73:VAL:HG11	1.59	0.83
49:DR:59:ILE:HG12	49:DR:101:ILE:HD13	1.60	0.83
50:DT:29:THR:HA	50:DT:86:THR:HA	1.58	0.83
1:AA:1432:G:H5'	28:BP:105:LYS:HG2	1.59	0.83
28:BP:56:SER:HB2	28:BP:75:THR:HB	1.60	0.83
3:CD:2:ARG:HH11	3:CD:114:ARG:HD3	1.41	0.83
1:CA:1115:U:H5'	9:CJ:68:ARG:HH22	1.44	0.83
29:BE:126:VAL:HG22	29:BE:127:GLU:H	1.44	0.83
40:BH:115:VAL:HB	40:BH:130:VAL:HG12	1.58	0.83
22:DA:32:U:H4'	22:DA:52:A:H62	1.42	0.83
47:DF:28:PRO:HB2	47:DF:168:LEU:HD21	1.60	0.83
8:AI:117:LEU:HD22	8:AI:123:ARG:HB3	1.61	0.83
40:BH:5:LEU:HD12	40:BH:17:ASP:HB3	1.58	0.83
50:BT:5:GLU:HA	50:BT:8:LEU:HB2	1.58	0.83
2:CC:106:ARG:NE	2:CC:106:ARG:H	1.76	0.83
2:AC:61:LYS:HB3	2:AC:96:VAL:HG11	1.59	0.83
1:AA:1533:C:O2'	1:AA:1534:A:H4'	1.79	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AE:132:PRO:HG2	4:AE:133:ILE:HD12	1.61	0.83
10:AK:108:ASN:HD21	19:AU:6:ARG:HD2	1.43	0.83
23:BB:1060:U:C2	23:BB:1088:A:N7	2.47	0.83
27:BK:54:LYS:H	27:BK:54:LYS:HD2	1.44	0.83
12:AM:90:HIS:HA	12:AM:108:ARG:HH22	1.44	0.82
40:BH:100:ALA:HB3	40:BH:112:LYS:HA	1.61	0.82
6:CG:120:ALA:HA	6:CG:123:LEU:HD12	1.58	0.82
23:BB:1082:U:C4	23:BB:1086:A:C2	2.66	0.82
23:BB:661:A:H1'	37:BL:12:SER:O	1.79	0.82
23:DB:2328:A:H2'	23:DB:2329:U:C6	2.14	0.82
29:DE:5:LEU:HD12	29:DE:10:SER:HB2	1.60	0.82
24:DI:27:LEU:HD23	24:DI:27:LEU:H	1.44	0.82
22:BA:89:U:O2	23:BB:958:U:H2'	1.79	0.82
29:BE:18:THR:HA	29:BE:106:LYS:HG3	1.62	0.82
29:BE:61:ARG:NH1	29:BE:64:GLY:HA3	1.93	0.82
51:BZ:7:VAL:HG13	51:BZ:8:THR:HG23	1.61	0.82
41:DJ:17:VAL:HG23	41:DJ:137:PRO:HB2	1.58	0.82
9:CJ:52:LEU:H	9:CJ:52:LEU:HD12	1.43	0.82
23:DB:1019:U:H2'	23:DB:1020:A:C8	2.15	0.82
1:AA:1063:C:H42	1:AA:1193:G:H1	1.23	0.82
23:BB:2328:A:H2'	23:BB:2329:U:C6	2.14	0.82
23:BB:2529:G:H4'	48:BG:174:LYS:HG3	1.61	0.82
23:DB:1024:G:H3'	23:DB:1025:G:H5''	1.59	0.82
48:DG:153:PRO:HG2	48:DG:162:ARG:HB3	1.62	0.82
28:DP:20:ARG:HG3	28:DP:21:PRO:HD2	1.61	0.82
48:BG:43:LYS:HB2	48:BG:50:THR:HB	1.61	0.82
23:DB:321:U:H1'	29:DE:162:ARG:HH11	1.44	0.82
40:DH:116:ARG:HB3	40:DH:131:SER:H	1.44	0.82
52:DW:39:GLN:HG2	52:DW:40:ARG:N	1.92	0.82
23:BB:1024:G:H3'	23:BB:1025:G:H5''	1.62	0.82
28:BP:21:PRO:HG2	28:BP:91:VAL:HG21	1.59	0.82
23:DB:1082:U:C4	23:DB:1086:A:C2	2.67	0.82
23:BB:1579:A:H2'	23:BB:1580:A:C8	2.15	0.82
38:BM:35:ALA:HB2	38:BM:100:LYS:HB2	1.61	0.82
23:DB:2674:G:H4'	27:DK:30:ARG:HG3	1.61	0.82
1:AA:922:G:H4'	4:AE:24:VAL:HA	1.62	0.82
42:BN:2:ARG:HG2	42:BN:5:LYS:HB2	1.62	0.82
3:CD:29:THR:H	3:CD:33:ILE:HG22	1.44	0.82
11:CL:51:VAL:HG12	11:CL:52:CYS:H	1.45	0.82
23:DB:1060:U:C2	23:DB:1088:A:N7	2.48	0.82
24:DI:121:ILE:HD13	24:DI:121:ILE:H	1.45	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DR:58:VAL:HG22	49:DR:59:ILE:H	1.44	0.82
52:DW:45:HIS:HB2	52:DW:50:VAL:HG12	1.62	0.82
23:BB:2531:A:H5'	48:BG:156:TYR:CZ	2.14	0.81
12:CM:104:ASN:H	12:CM:106:ARG:HH21	1.28	0.81
23:DB:196:A:H5''	37:DL:47:ARG:HH12	1.43	0.81
12:AM:38:ILE:HG13	12:AM:55:LEU:HD21	1.60	0.81
49:BR:39:LEU:HA	49:BR:53:PHE:HA	1.60	0.81
23:DB:448:U:H5	23:DB:583:G:H1'	1.45	0.81
29:DE:18:THR:HA	29:DE:106:LYS:HG3	1.61	0.81
37:DL:124:GLY:N	37:DL:143:GLU:HG3	1.94	0.81
43:DO:47:VAL:HG12	43:DO:48:LEU:H	1.45	0.81
8:AI:113:LYS:HG2	8:AI:120:ALA:H	1.46	0.81
1:AA:1289:A:H61	8:AI:71:ILE:HD11	1.45	0.81
10:AK:33:ILE:HB	10:AK:73:VAL:HG11	1.61	0.81
37:BL:29:LYS:HG3	37:BL:30:THR:HG23	1.62	0.81
13:CP:5:ARG:NH2	13:CP:24:SER:HA	1.95	0.81
5:AF:11:HIS:HB3	5:AF:14:GLN:HG3	1.61	0.81
7:AH:86:LYS:HB3	7:AH:90:GLU:HB3	1.60	0.81
32:B4:2:LYS:HG2	32:B4:4:ARG:HE	1.44	0.81
23:BB:2010:G:H5''	45:BS:42:LYS:HB2	1.63	0.81
35:BV:9:ARG:HH21	35:BV:12:GLN:HA	1.45	0.81
17:AT:34:VAL:HG11	17:AT:78:LEU:HD22	1.63	0.81
40:DH:5:LEU:HD12	40:DH:17:ASP:HB3	1.59	0.81
27:BK:105:ARG:HD3	27:BK:105:ARG:H	1.42	0.81
23:BB:571:U:H3'	49:BR:80:ARG:HH12	1.46	0.81
1:CA:1288:A:N1	1:CA:1371:G:H1'	1.96	0.81
1:CA:1492:A:H3'	1:CA:1493:A:O4'	1.80	0.81
23:DB:571:U:H3'	49:DR:80:ARG:HH12	1.44	0.81
1:CA:975:A:H4'	1:CA:976:G:O5'	1.78	0.81
29:DE:111:GLU:HA	29:DE:114:ARG:HG3	1.62	0.81
29:DE:58:LYS:H	29:DE:58:LYS:NZ	1.78	0.81
47:DF:84:ILE:HG23	47:DF:85:GLY:H	1.45	0.81
48:DG:8:VAL:HG11	48:DG:49:LEU:HB3	1.61	0.81
12:AM:33:LEU:HD22	12:AM:38:ILE:HB	1.61	0.81
47:BF:28:PRO:HB2	47:BF:168:LEU:HD21	1.63	0.81
48:BG:153:PRO:HG2	48:BG:162:ARG:HB3	1.63	0.81
12:CM:14:ALA:HB2	12:CM:42:VAL:HG23	1.63	0.81
29:DE:4:VAL:HG12	29:DE:5:LEU:H	1.46	0.81
48:DG:148:ARG:HD3	48:DG:152:ARG:HH11	1.45	0.81
48:DG:43:LYS:HB2	48:DG:50:THR:HB	1.61	0.81
1:AA:279:A:H5''	1:AA:280:C:H3'	1.61	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:29:VAL:HB	26:BD:98:VAL:HG22	1.63	0.81
17:CT:75:LYS:HA	17:CT:78:LEU:HD12	1.61	0.81
9:AJ:17:LEU:HD12	9:AJ:96:VAL:HG13	1.61	0.81
36:B2:45:SER:O	36:B2:46:LYS:HB2	1.80	0.81
23:BB:141:G:H5''	23:BB:142:A:O4'	1.82	0.81
47:BF:137:PHE:HB2	47:BF:138:PRO:HD2	1.62	0.81
45:BS:26:GLY:H	45:BS:71:VAL:HG13	1.43	0.81
1:CA:21:G:H2'	1:CA:22:G:C8	2.16	0.81
1:CA:958:A:H61	16:CS:76:THR:HG22	1.44	0.81
36:D2:45:SER:O	36:D2:46:LYS:HB2	1.80	0.81
29:DE:3:LEU:H	29:DE:13:THR:H	1.29	0.81
42:DN:12:ARG:HG2	42:DN:16:HIS:ND1	1.96	0.81
4:AE:28:ARG:NH2	4:AE:30:PHE:HA	1.96	0.80
40:BH:31:VAL:HB	40:BH:32:PRO:CD	2.11	0.80
23:DB:674:G:H1'	29:DE:69:ARG:HE	1.45	0.80
3:AD:2:ARG:HH11	3:AD:114:ARG:HD3	1.44	0.80
23:BB:2133:G:N3	23:BB:2133:G:H2'	1.95	0.80
23:BB:543:G:H2'	23:BB:544:C:H5''	1.63	0.80
12:CM:21:ILE:HG22	12:CM:23:GLY:H	1.44	0.80
18:AB:57:ASN:HB3	18:AB:219:THR:HG22	1.63	0.80
11:AL:51:VAL:HG12	11:AL:52:CYS:H	1.45	0.80
12:AM:49:GLU:O	12:AM:52:ILE:HG22	1.80	0.80
29:BE:4:VAL:HG12	29:BE:5:LEU:H	1.45	0.80
40:BH:6:LEU:HD11	40:BH:37:VAL:HG12	1.62	0.80
43:BO:39:VAL:HG12	43:BO:48:LEU:HD12	1.63	0.80
1:CA:676:A:H1'	10:CK:116:PRO:HB3	1.62	0.80
3:CD:160:LEU:H	3:CD:160:LEU:HD13	1.46	0.80
27:DK:71:ARG:HB3	27:DK:72:PRO:HD2	1.61	0.80
4:AE:71:ILE:HD11	4:AE:144:GLU:HG3	1.63	0.80
29:BE:149:ILE:HD11	29:BE:172:ALA:HA	1.62	0.80
27:BK:38:ILE:H	27:BK:38:ILE:HD13	1.46	0.80
4:CE:28:ARG:NH2	4:CE:30:PHE:HA	1.96	0.80
8:CI:16:ALA:HB2	8:CI:66:VAL:HG23	1.62	0.80
10:CK:105:ARG:HH21	19:CU:10:PRO:HB3	1.45	0.80
23:DB:2722:G:H4'	42:DN:4:ARG:HB2	1.63	0.80
40:DH:31:VAL:HB	40:DH:32:PRO:CD	2.11	0.80
12:AM:44:ILE:HA	12:AM:47:LEU:HD23	1.64	0.80
23:BB:718:A:H3'	23:BB:719:C:H6	1.45	0.80
49:BR:58:VAL:HG22	49:BR:59:ILE:H	1.46	0.80
27:DK:107:LEU:HD12	27:DK:107:LEU:H	1.47	0.80
1:AA:17:U:H2'	1:AA:18:C:C6	2.16	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:125:PHE:HD2	18:AB:125:PHE:H	1.29	0.80
13:AP:5:ARG:NH2	13:AP:24:SER:HA	1.96	0.80
23:BB:1652:A:H62	42:BN:11:ASN:HD21	1.28	0.80
23:DB:1099:G:H8	24:DI:3:LYS:CA	1.95	0.80
23:DB:1168:G:H2'	23:DB:1169:A:C8	2.17	0.80
23:DB:876:C:H42	23:DB:901:C:N4	1.80	0.80
24:DI:21:PRO:HB2	24:DI:22:PRO:HD3	1.63	0.80
6:AG:145:GLU:HA	6:AG:148:LYS:HB2	1.64	0.80
21:CN:65:GLN:NE2	21:CN:65:GLN:H	1.79	0.80
47:DF:90:LEU:HD12	47:DF:94:ARG:HB3	1.62	0.80
9:AJ:7:ARG:HB2	9:AJ:101:SER:N	1.94	0.80
23:BB:2799:A:H4'	23:BB:2800:A:C8	2.17	0.80
40:BH:73:ASN:N	40:BH:73:ASN:HD22	1.78	0.80
23:DB:2472:G:H2'	23:DB:2475:C:H42	1.45	0.80
35:DV:9:ARG:HH21	35:DV:12:GLN:HA	1.47	0.80
23:BB:2472:G:H2'	23:BB:2475:C:H42	1.46	0.80
23:BB:321:U:H1'	29:BE:162:ARG:HH11	1.47	0.80
6:CG:19:SER:HB3	6:CG:22:LEU:HB2	1.62	0.80
44:BQ:8:ILE:H	44:BQ:8:ILE:HD12	1.45	0.80
1:CA:1313:U:OP2	16:CS:5:LYS:HA	1.82	0.80
1:CA:1499:A:H2'	1:CA:1500:A:H8	1.47	0.80
22:DA:89:U:O2	23:DB:958:U:H2'	1.82	0.80
47:DF:137:PHE:HB2	47:DF:138:PRO:HD2	1.62	0.80
48:BG:8:VAL:HG11	48:BG:49:LEU:HB3	1.64	0.79
52:DW:48:ALA:HB3	52:DW:81:ILE:HG13	1.63	0.79
1:AA:1221:G:H4'	16:AS:76:THR:HG21	1.63	0.79
1:CA:1206:G:H4'	2:CC:192:TYR:HA	1.64	0.79
9:CJ:92:LEU:H	9:CJ:92:LEU:HD22	1.46	0.79
23:BB:670:A:H4'	23:BB:671:C:H5'	1.64	0.79
52:BW:48:ALA:HB3	52:BW:81:ILE:HG13	1.65	0.79
5:CF:11:HIS:HB3	5:CF:14:GLN:HG3	1.64	0.79
48:DG:6:ALA:HB3	48:DG:68:ARG:HG3	1.65	0.79
3:AD:160:LEU:H	3:AD:160:LEU:HD13	1.45	0.79
8:AI:12:LYS:H	8:AI:105:ARG:NH2	1.80	0.79
11:AL:24:GLU:HB2	11:AL:26:CYS:SG	2.22	0.79
17:AT:75:LYS:HA	17:AT:78:LEU:HD12	1.64	0.79
37:BL:124:GLY:N	37:BL:143:GLU:HG3	1.98	0.79
1:CA:13:U:H1'	1:CA:914:A:H5''	1.64	0.79
21:CN:30:ILE:HG22	21:CN:44:VAL:HG21	1.64	0.79
29:DE:117:ARG:HA	29:DE:185:LYS:HG2	1.65	0.79
11:AL:78:VAL:HG12	11:AL:101:LEU:HD13	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BG:6:ALA:HB3	48:BG:68:ARG:HG3	1.64	0.79
22:BA:11:C:H5'	52:BW:71:LYS:HE3	1.64	0.79
51:BZ:4:VAL:HG12	51:BZ:11:ARG:HG2	1.62	0.79
21:CN:40:ARG:HH11	21:CN:40:ARG:HB2	1.47	0.79
32:D4:2:LYS:HG2	32:D4:4:ARG:HE	1.48	0.79
1:AA:21:G:H2'	1:AA:22:G:C8	2.17	0.79
2:AC:6:PRO:HG3	2:AC:200:TRP:HE1	1.48	0.79
9:AJ:26:VAL:HG12	9:AJ:30:LYS:HE2	1.63	0.79
9:AJ:42:LEU:HD23	9:AJ:71:LEU:HD11	1.64	0.79
48:BG:84:LYS:HG2	48:BG:85:LYS:H	1.47	0.79
47:DF:110:ILE:HA	47:DF:111:ARG:CZ	2.12	0.79
44:DQ:57:ARG:HH11	44:DQ:61:ILE:HD11	1.47	0.79
54:AA:1661:SCM:H2M1	4:AE:26:GLY:HA2	1.63	0.79
3:AD:29:THR:H	3:AD:33:ILE:HG22	1.47	0.79
10:AK:91:GLY:HA2	10:AK:94:SER:HB3	1.65	0.79
48:DG:84:LYS:HG2	48:DG:85:LYS:H	1.48	0.79
5:AF:38:ARG:HB3	5:AF:63:ASN:HB2	1.65	0.79
41:BJ:81:ILE:HG23	41:BJ:82:GLY:H	1.48	0.79
41:DJ:81:ILE:HG23	41:DJ:82:GLY:H	1.47	0.79
47:BF:2:LYS:HD2	47:BF:100:GLU:HG2	1.63	0.79
47:BF:90:LEU:HD12	47:BF:94:ARG:HB3	1.64	0.79
24:BI:21:PRO:HB2	24:BI:22:PRO:HD3	1.62	0.79
4:CE:71:ILE:HD11	4:CE:144:GLU:HG3	1.65	0.79
16:AS:66:VAL:HG23	16:AS:67:GLY:H	1.47	0.79
42:BN:12:ARG:HG2	42:BN:16:HIS:ND1	1.98	0.79
1:CA:337:G:H2'	1:CA:338:A:H8	1.48	0.79
1:CA:462:G:H5'	1:CA:463:U:OP2	1.83	0.79
2:CC:120:THR:HG22	2:CC:188:ALA:HB2	1.65	0.79
11:CL:29:LYS:HB3	11:CL:56:LEU:HD22	1.64	0.79
5:AF:3:HIS:HB2	5:AF:92:THR:HA	1.64	0.78
23:DB:670:A:H4'	23:DB:671:C:H5'	1.65	0.78
23:DB:1099:G:P	24:DI:4:VAL:H	2.06	0.78
1:AA:13:U:H1'	1:AA:914:A:H5''	1.64	0.78
23:BB:2376:A:H61	43:BO:94:ARG:HD2	1.48	0.78
15:CR:38:ILE:H	15:CR:38:ILE:HD13	1.48	0.78
25:DC:172:THR:HG22	25:DC:182:LYS:HG2	1.64	0.78
40:DH:77:THR:HG22	40:DH:143:ILE:HB	1.66	0.78
1:AA:676:A:H1'	10:AK:116:PRO:HB3	1.63	0.78
23:BB:1019:U:H2'	23:BB:1020:A:C8	2.18	0.78
25:BC:172:THR:HG22	25:BC:182:LYS:HG2	1.64	0.78
26:BD:34:VAL:CG1	26:BD:94:GLN:H	1.96	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1292:G:H2'	1:CA:1293:C:C6	2.19	0.78
8:CI:29:ILE:HG12	8:CI:64:ILE:HB	1.65	0.78
9:CJ:8:ILE:HA	9:CJ:100:ILE:HG22	1.64	0.78
13:AP:4:ILE:HG12	13:AP:21:VAL:HG22	1.65	0.78
18:CB:216:VAL:O	18:CB:220:VAL:HG23	1.84	0.78
26:DD:106:LYS:HB3	26:DD:206:ALA:H	1.48	0.78
27:DK:38:ILE:HD13	27:DK:38:ILE:H	1.47	0.78
1:AA:764:C:H3'	1:AA:765:G:H21	1.48	0.78
18:AB:163:ILE:HG23	18:AB:164:ASP:H	1.49	0.78
2:AC:126:ARG:HA	2:AC:126:ARG:HH11	1.49	0.78
21:AN:31:SER:H	21:AN:44:VAL:HG11	1.48	0.78
23:BB:2060:A:H2'	29:BE:63:LYS:HE2	1.63	0.78
24:BI:33:ASN:HD21	24:BI:64:ARG:HH11	1.31	0.78
27:BK:19:VAL:HG13	27:BK:43:ILE:HA	1.66	0.78
27:BK:71:ARG:HB3	27:BK:72:PRO:HD2	1.64	0.78
5:CF:38:ARG:HB3	5:CF:63:ASN:HB2	1.65	0.78
32:D4:36:ARG:HG2	32:D4:37:GLN:H	1.48	0.78
44:DQ:8:ILE:HD12	44:DQ:8:ILE:H	1.46	0.78
1:AA:1101:A:H4'	1:AA:1102:A:O5'	1.84	0.78
40:BH:90:LEU:HD13	40:BH:123:ARG:HA	1.65	0.78
18:AB:120:SER:HA	18:AB:125:PHE:CB	2.14	0.78
8:AI:118:ARG:HB2	8:AI:124:PRO:HD3	1.64	0.78
23:BB:2306:C:H3'	23:BB:2307:G:C5'	2.13	0.78
44:BQ:30:VAL:HG22	44:BQ:31:TYR:H	1.48	0.78
2:CC:113:LYS:HD2	2:CC:184:ASN:HD21	1.46	0.78
25:DC:144:GLU:HG3	25:DC:151:GLY:N	1.98	0.78
48:DG:26:LYS:HA	48:DG:32:LEU:H	1.48	0.78
43:DO:39:VAL:HG12	43:DO:48:LEU:HD12	1.65	0.78
32:B4:36:ARG:HG2	32:B4:37:GLN:H	1.49	0.78
48:BG:15:ASP:HB2	48:BG:26:LYS:HD3	1.66	0.78
24:BI:55:PRO:HD3	24:BI:74:PRO:HD3	1.64	0.78
37:BL:6:LEU:H	37:BL:6:LEU:HD23	1.48	0.78
28:BP:20:ARG:HG3	28:BP:21:PRO:HD2	1.65	0.78
12:CM:54:THR:O	12:CM:58:GLU:HB2	1.83	0.78
17:CT:34:VAL:HG11	17:CT:78:LEU:HD22	1.64	0.78
1:AA:376:G:H5''	13:AP:5:ARG:HB2	1.66	0.78
6:AG:13:PRO:HA	6:AG:20:GLU:HA	1.65	0.78
6:AG:71:THR:H	6:AG:141:HIS:HE1	1.30	0.78
16:AS:29:PRO:HB3	16:AS:47:THR:HB	1.64	0.78
1:CA:1004:A:H5'	1:CA:1024:G:H22	1.47	0.78
1:CA:1483:A:H2'	1:CA:1484:C:O4'	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:196:ASN:ND2	25:DC:199:HIS:HB2	1.99	0.78
1:CA:1238:A:H5'	1:CA:1336:C:N4	1.99	0.78
18:CB:14:HIS:HD2	18:CB:202:ASN:H	1.30	0.78
10:CK:108:ASN:HD21	19:CU:6:ARG:HD2	1.48	0.78
44:DQ:89:ILE:HG21	49:DR:11:GLN:HE22	1.47	0.78
51:DZ:40:VAL:HG21	51:DZ:43:GLU:HB3	1.65	0.78
1:AA:68:G:H5'	1:AA:171:A:H1'	1.66	0.77
1:CA:1236:A:H4'	1:CA:1304:G:H4'	1.66	0.77
19:AU:11:PHE:HD2	2:CC:108:PRO:HD2	1.49	0.77
26:DD:29:VAL:HB	26:DD:98:VAL:HG22	1.64	0.77
37:BL:57:LEU:HD12	37:BL:60:ARG:HH22	1.49	0.77
1:CA:83:C:O2'	1:CA:84:U:H2'	1.84	0.77
47:DF:62:GLN:HG3	47:DF:91:ARG:HH11	1.49	0.77
39:DX:3:ALA:HA	39:DX:6:LEU:HD23	1.64	0.77
23:BB:979:A:H2'	23:BB:982:C:H41	1.50	0.77
1:CA:979:C:H1'	1:CA:1317:C:H41	1.48	0.77
23:DB:2306:C:H3'	23:DB:2307:G:C5'	2.13	0.77
23:DB:2742:G:OP1	32:D4:36:ARG:HD3	1.81	0.77
40:DH:4:ILE:HB	40:DH:37:VAL:HG12	1.67	0.77
18:AB:166:ASP:OD1	18:AB:190:SER:HA	1.85	0.77
18:AB:86:CYS:HB3	18:AB:88:GLN:NE2	1.99	0.77
23:BB:590:A:H2'	23:BB:591:U:C6	2.20	0.77
48:BG:132:LEU:HD23	48:BG:132:LEU:H	1.48	0.77
23:DB:1099:G:H5''	24:DI:2:LYS:C	2.04	0.77
23:DB:1558:C:H4'	23:DB:1559:U:H5'	1.65	0.77
23:DB:947:A:H2'	23:DB:948:C:C6	2.19	0.77
23:BB:1387:A:H5'	23:BB:1469:A:H1'	1.67	0.77
26:BD:118:PHE:CE2	42:BN:1:MET:HB3	2.20	0.77
37:BL:80:SER:HA	37:BL:115:GLU:HB2	1.67	0.77
1:CA:1342:C:O2'	8:CI:125:GLN:HB3	1.84	0.77
8:CI:50:PRO:HD3	8:CI:79:ARG:HG3	1.66	0.77
23:DB:1199:U:H5'	44:DQ:4:LYS:HD3	1.67	0.77
44:DQ:30:VAL:HG22	44:DQ:31:TYR:H	1.49	0.77
23:BB:1060:U:OP2	24:BI:74:PRO:HA	1.84	0.77
45:BS:24:ILE:HG22	45:BS:71:VAL:HG11	1.65	0.77
50:BT:73:ARG:HH21	50:BT:73:ARG:HB3	1.49	0.77
46:BU:47:PRO:HB2	46:BU:53:GLN:HB2	1.67	0.77
25:DC:2:VAL:HG23	25:DC:3:VAL:H	1.50	0.77
29:DE:61:ARG:NH1	29:DE:64:GLY:HA3	2.00	0.77
45:DS:24:ILE:HG22	45:DS:71:VAL:HG11	1.66	0.77
25:BC:94:LEU:HD13	25:BC:100:ARG:HD2	1.64	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BO:47:VAL:HG12	43:BO:48:LEU:H	1.50	0.77
52:BW:51:GLY:CA	52:BW:59:PHE:HB2	2.15	0.77
12:CM:29:SER:HA	12:CM:32:ILE:HG22	1.65	0.77
23:DB:899:A:H3'	23:DB:900:A:H8	1.48	0.77
23:DB:600:G:H1'	29:DE:100:MET:HG2	1.65	0.77
24:DI:55:PRO:HD3	24:DI:74:PRO:HD3	1.66	0.77
1:AA:203:G:H1'	1:AA:465:A:C6	2.20	0.77
8:AI:55:ASP:HB2	8:AI:59:LYS:HD2	1.64	0.77
14:AQ:20:ILE:HD13	14:AQ:47:ASP:HB3	1.67	0.77
33:D1:34:GLU:HG2	33:D1:49:LYS:HG3	1.67	0.77
18:AB:23:ASN:HD21	18:AB:25:LYS:HB2	1.49	0.77
3:AD:90:LEU:HA	3:AD:93:LEU:HD12	1.66	0.77
12:AM:78:ARG:O	12:AM:82:LEU:HG	1.84	0.77
23:BB:1558:C:H4'	23:BB:1559:U:H5'	1.65	0.77
23:BB:2901:C:H2'	23:BB:2902:C:H5'	1.67	0.77
52:BW:45:HIS:HB2	52:BW:50:VAL:HG12	1.65	0.77
1:CA:176:C:H2'	1:CA:177:G:N3	2.00	0.77
1:CA:243:A:H4'	1:CA:244:U:H5'	1.67	0.77
20:CO:71:LYS:HG2	20:CO:75:VAL:HG22	1.67	0.77
23:DB:1097:U:H2'	23:DB:1098:A:O4'	1.84	0.77
23:DB:571:U:H3'	49:DR:80:ARG:NH1	2.00	0.77
26:DD:151:THR:HB	26:DD:152:PRO:HD3	1.67	0.77
29:DE:2:GLU:HA	29:DE:13:THR:HA	1.66	0.77
18:AB:13:VAL:HB	18:AB:207:ARG:HE	1.50	0.77
18:AB:61:SER:HB2	18:AB:62:ARG:NH1	2.00	0.77
23:BB:448:U:H5	23:BB:583:G:H1'	1.50	0.77
26:BD:106:LYS:HB3	26:BD:206:ALA:H	1.50	0.77
8:CI:19:PHE:HB2	8:CI:63:TYR:HB3	1.66	0.77
8:CI:51:LEU:HD13	8:CI:56:MET:HG2	1.67	0.77
23:DB:1387:A:H5'	23:DB:1469:A:H1'	1.66	0.77
24:DI:45:THR:HA	24:DI:48:ILE:HG22	1.66	0.77
49:DR:72:VAL:HG23	49:DR:89:HIS:HB3	1.67	0.77
2:AC:123:LEU:HA	2:AC:127:VAL:CG2	2.14	0.76
29:BE:147:LEU:HB3	29:BE:186:VAL:HG23	1.66	0.76
47:BF:8:LYS:HA	47:BF:12:VAL:HG21	1.67	0.76
45:BS:4:ILE:HG22	45:BS:106:VAL:HA	1.65	0.76
10:CK:22:ILE:HB	10:CK:85:VAL:HG22	1.67	0.76
23:DB:1373:A:H2'	23:DB:1374:G:O4'	1.84	0.76
23:DB:718:A:H3'	23:DB:719:C:H6	1.48	0.76
37:DL:6:LEU:H	37:DL:6:LEU:HD23	1.50	0.76
48:BG:96:ALA:HB3	48:BG:103:ASN:HB3	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2880:C:O4'	42:BN:91:ALA:HB3	1.85	0.76
44:BQ:57:ARG:HH11	44:BQ:61:ILE:HD11	1.51	0.76
30:BY:6:ILE:HD13	30:BY:6:ILE:H	1.50	0.76
23:DB:1099:G:OP2	24:DI:3:LYS:HA	1.84	0.76
52:DW:51:GLY:CA	52:DW:59:PHE:HB2	2.14	0.76
1:AA:1499:A:H2'	1:AA:1500:A:H8	1.50	0.76
29:BE:117:ARG:HA	29:BE:185:LYS:HG2	1.65	0.76
29:BE:3:LEU:H	29:BE:13:THR:H	1.28	0.76
46:BU:85:ARG:HA	46:BU:85:ARG:NE	2.00	0.76
34:D3:31:ILE:HD11	34:D3:34:LYS:HD3	1.67	0.76
32:D4:5:ALA:HA	32:D4:37:GLN:HE22	1.49	0.76
23:DB:979:A:H2'	23:DB:982:C:H41	1.50	0.76
23:DB:2330:G:H1'	52:DW:38:ARG:HB2	1.68	0.76
33:B1:5:ARG:HH11	33:B1:25:ASN:HB2	1.49	0.76
32:B4:5:ALA:HA	32:B4:37:GLN:HE22	1.49	0.76
23:BB:2039:U:H2'	23:BB:2040:G:H8	1.51	0.76
25:BC:2:VAL:HG23	25:BC:3:VAL:H	1.51	0.76
47:BF:110:ILE:HA	47:BF:111:ARG:CZ	2.15	0.76
42:BN:2:ARG:HA	42:BN:5:LYS:HD3	1.65	0.76
46:BU:82:VAL:H	46:BU:96:LYS:HZ2	1.34	0.76
52:BW:27:GLY:O	52:BW:63:ASP:HA	1.85	0.76
1:CA:67:C:H2'	1:CA:68:G:H8	1.50	0.76
18:CB:48:MET:O	18:CB:51:GLU:HB3	1.85	0.76
3:CD:26:ALA:HA	3:CD:30:LYS:HE3	1.67	0.76
10:CK:91:GLY:HA2	10:CK:94:SER:HB3	1.68	0.76
26:DD:34:VAL:CG1	26:DD:94:GLN:H	1.99	0.76
23:DB:1099:G:P	24:DI:3:LYS:HA	2.25	0.76
1:AA:269:C:H2'	1:AA:270:A:C8	2.20	0.76
18:AB:107:ARG:HH22	18:AB:111:LYS:HD3	1.49	0.76
33:B1:37:LYS:H	33:B1:48:TYR:HD2	1.34	0.76
39:BX:3:ALA:HA	39:BX:6:LEU:HD23	1.66	0.76
1:CA:67:C:H2'	1:CA:68:G:C8	2.20	0.76
16:CS:10:ILE:HG22	16:CS:14:LEU:HD21	1.67	0.76
29:DE:3:LEU:HB2	29:DE:12:LEU:HD12	1.68	0.76
52:DW:27:GLY:O	52:DW:63:ASP:HA	1.84	0.76
8:AI:51:LEU:HB3	8:AI:56:MET:HB3	1.65	0.76
3:CD:12:ARG:HH11	3:CD:12:ARG:HB3	1.51	0.76
8:CI:18:VAL:HG11	8:CI:82:ILE:HA	1.66	0.76
8:CI:66:VAL:HG21	8:CI:74:GLN:HG3	1.66	0.76
29:DE:149:ILE:HD11	29:DE:172:ALA:HA	1.65	0.76
20:AO:32:LEU:O	20:AO:36:ILE:HG12	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AS:17:LYS:HD2	16:AS:30:LEU:HD22	1.66	0.76
23:BB:1046:A:C3'	23:BB:1047:G:H5''	2.15	0.76
23:BB:1723:G:H3'	23:BB:1724:G:H8	1.50	0.76
23:BB:79:C:O2'	23:BB:346:A:H1'	1.85	0.76
45:DS:4:ILE:HG22	45:DS:106:VAL:HA	1.66	0.76
1:AA:1213:A:H2'	1:AA:1215:G:N7	2.00	0.76
15:AR:38:ILE:H	15:AR:38:ILE:HD13	1.50	0.76
23:BB:1082:U:N3	23:BB:1086:A:C2	2.53	0.76
47:BF:147:ARG:HG2	47:BF:148:VAL:HG22	1.68	0.76
40:BH:99:ILE:HG13	40:BH:130:VAL:HG21	1.67	0.76
51:BZ:5:CYS:HB2	51:BZ:10:LYS:HB2	1.66	0.76
12:CM:12:LYS:HB3	12:CM:16:ILE:HG23	1.67	0.76
21:CN:41:TRP:HA	21:CN:41:TRP:CE3	2.19	0.76
20:CO:64:ARG:HE	20:CO:64:ARG:HA	1.50	0.76
16:CS:62:THR:HG22	16:CS:63:ASP:H	1.49	0.76
33:D1:33:LEU:HB3	33:D1:51:ALA:CB	2.16	0.76
23:DB:2039:U:H2'	23:DB:2040:G:H8	1.50	0.76
23:DB:2886:A:H62	31:D0:39:ARG:NE	1.83	0.76
44:DQ:26:ALA:O	44:DQ:30:VAL:HG12	1.86	0.76
8:AI:9:GLY:HA2	8:AI:80:HIS:HB3	1.68	0.76
23:BB:580:U:H2'	23:BB:581:C:C6	2.20	0.76
29:BE:148:ILE:HG13	29:BE:167:VAL:HG23	1.68	0.76
1:CA:764:C:H3'	1:CA:765:G:H21	1.50	0.76
3:CD:90:LEU:HA	3:CD:93:LEU:HD12	1.68	0.76
12:CM:28:ARG:NH1	12:CM:32:ILE:HD12	2.00	0.76
23:DB:919:U:H2'	23:DB:920:A:C8	2.20	0.76
46:DU:85:ARG:NE	46:DU:85:ARG:HA	2.01	0.76
30:DY:6:ILE:HD13	30:DY:6:ILE:H	1.49	0.76
1:AA:1391:U:H2'	1:AA:1392:G:C8	2.19	0.76
34:B3:31:ILE:HD11	34:B3:34:LYS:HD3	1.66	0.76
23:BB:2134:A:N6	23:BB:2157:G:H21	1.83	0.76
26:BD:116:LYS:HG3	26:BD:165:MET:SD	2.26	0.76
48:BG:26:LYS:HA	48:BG:32:LEU:H	1.50	0.76
1:CA:337:G:H2'	1:CA:338:A:C8	2.20	0.76
10:CK:28:ASN:ND2	10:CK:46:ALA:HB3	2.00	0.76
23:DB:2267:A:C8	23:DB:2267:A:H3'	2.21	0.76
48:DG:9:VAL:HA	48:DG:48:THR:HB	1.67	0.76
6:AG:47:GLU:HA	6:AG:57:GLU:HG2	1.68	0.75
12:AM:21:ILE:HG23	12:AM:65:GLU:OE2	1.86	0.75
14:AQ:13:SER:HB3	14:AQ:21:VAL:HB	1.65	0.75
23:BB:947:A:H2'	23:BB:948:C:C6	2.21	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CB:212:TYR:HA	18:CB:215:ALA:HB3	1.66	0.75
23:DB:1251:C:H2'	44:DQ:5:ARG:HH12	1.51	0.75
25:DC:4:LYS:HD2	25:DC:5:CYS:H	1.50	0.75
26:DD:53:GLY:HA3	26:DD:77:ARG:HG3	1.68	0.75
1:AA:946:A:H2'	1:AA:947:G:C8	2.20	0.75
1:AA:1348:U:H4'	8:AI:121:ARG:HG3	1.65	0.75
10:AK:19:VAL:HG23	10:AK:34:THR:HG23	1.68	0.75
20:AO:22:THR:HA	20:AO:27:VAL:HG11	1.68	0.75
13:AP:28:ARG:HD2	13:AP:29:ASN:N	2.00	0.75
40:BH:72:ILE:HG13	40:BH:75:LEU:HD21	1.68	0.75
1:CA:9:G:H5'	4:CE:107:GLY:HA3	1.68	0.75
18:CB:100:LEU:HG	18:CB:178:LEU:HD11	1.68	0.75
40:DH:6:LEU:HD11	40:DH:37:VAL:HG12	1.67	0.75
24:DI:9:LYS:HG2	24:DI:57:VAL:HG13	1.69	0.75
50:DT:73:ARG:HH21	50:DT:73:ARG:HB3	1.50	0.75
9:AJ:39:PRO:HA	9:AJ:74:VAL:HA	1.67	0.75
29:BE:63:LYS:HA	29:BE:63:LYS:HE3	1.68	0.75
1:CA:376:G:H5''	13:CP:5:ARG:HB2	1.66	0.75
16:CS:10:ILE:HG12	16:CS:40:PHE:CZ	2.20	0.75
23:DB:2008:C:H2'	23:DB:2009:A:H8	1.51	0.75
26:DD:9:VAL:HA	26:DD:197:THR:HG23	1.65	0.75
29:DE:147:LEU:HB3	29:DE:186:VAL:HG23	1.66	0.75
47:DF:2:LYS:HD2	47:DF:100:GLU:HG2	1.67	0.75
35:DV:44:HIS:HE1	35:DV:86:LEU:H	1.34	0.75
6:AG:138:GLU:HB3	6:AG:142:ARG:HH21	1.51	0.75
8:AI:11:ARG:HG2	8:AI:77:ALA:HB2	1.68	0.75
23:BB:2886:A:H62	31:B0:39:ARG:NE	1.83	0.75
11:CL:14:LYS:HG2	11:CL:16:ALA:H	1.51	0.75
23:DB:1141:U:H4'	23:DB:1142:A:O4'	1.85	0.75
48:DG:15:ASP:HB2	48:DG:26:LYS:HD3	1.68	0.75
1:AA:239:U:H4'	1:AA:239:U:OP1	1.86	0.75
25:BC:233:GLY:H	25:BC:241:LYS:HZ2	1.33	0.75
26:BD:53:GLY:HA3	26:BD:77:ARG:HG3	1.69	0.75
18:CB:9:LEU:HD23	18:CB:10:LYS:HG3	1.69	0.75
19:CU:34:ARG:HE	19:CU:35:GLU:C	1.89	0.75
40:DH:135:HIS:HB3	40:DH:138:VAL:HB	1.67	0.75
23:BB:1203:U:H1'	37:BL:4:ASN:HD21	1.50	0.75
47:BF:62:GLN:HG3	47:BF:91:ARG:HH11	1.50	0.75
28:BP:4:ILE:HG22	28:BP:5:LYS:H	1.51	0.75
44:BQ:26:ALA:O	44:BQ:30:VAL:HG12	1.87	0.75
45:BS:24:ILE:HG23	45:BS:32:ALA:HB1	1.69	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:269:C:H2'	1:CA:270:A:C8	2.21	0.75
2:CC:105:VAL:HA	2:CC:106:ARG:NH2	2.02	0.75
12:CM:47:LEU:HD13	12:CM:51:GLN:HB2	1.68	0.75
46:DU:47:PRO:HB2	46:DU:53:GLN:HB2	1.68	0.75
51:DZ:5:CYS:HB2	51:DZ:10:LYS:HB2	1.67	0.75
18:AB:112:ARG:HA	18:AB:115:ASP:OD2	1.87	0.75
11:AL:29:LYS:HB3	11:AL:56:LEU:HD22	1.68	0.75
19:AU:34:ARG:HE	19:AU:35:GLU:C	1.89	0.75
23:BB:1373:A:H2'	23:BB:1374:G:O4'	1.87	0.75
23:BB:2092:U:H4'	23:BB:2093:G:O5'	1.86	0.75
1:CA:301:G:H2'	1:CA:302:G:H8	1.51	0.75
2:CC:106:ARG:HD2	2:CC:106:ARG:O	1.87	0.75
50:BT:69:ARG:HE	50:BT:69:ARG:HA	1.52	0.75
21:CN:41:TRP:HA	21:CN:41:TRP:HE3	1.51	0.75
23:DB:1723:G:H3'	23:DB:1724:G:H8	1.52	0.75
23:DB:2787:C:H5'	26:DD:66:GLY:HA3	1.68	0.75
44:DQ:90:ASP:O	44:DQ:94:LEU:HB2	1.87	0.75
1:AA:1225:A:H5''	1:AA:1226:C:H5	1.51	0.75
1:AA:462:G:H5'	1:AA:463:U:OP2	1.87	0.75
1:AA:927:G:H4'	1:AA:1503:A:N7	2.02	0.75
8:AI:19:PHE:HB2	8:AI:63:TYR:HB3	1.67	0.75
23:BB:1341:G:H5'	50:BT:61:LEU:HD13	1.69	0.75
35:BV:72:VAL:HG21	35:BV:91:PHE:HB3	1.68	0.75
2:CC:13:ILE:H	2:CC:13:ILE:HD13	1.52	0.75
23:DB:358:U:H2'	23:DB:359:G:C8	2.22	0.75
28:DP:4:ILE:HG22	28:DP:5:LYS:H	1.52	0.75
1:AA:373:A:H1'	1:AA:481:G:N3	2.02	0.74
2:AC:190:THR:HG22	2:AC:191:THR:H	1.50	0.74
7:AH:94:VAL:HG23	7:AH:101:ALA:HB2	1.69	0.74
23:BB:919:U:H2'	23:BB:920:A:C8	2.21	0.74
1:CA:1317:C:H2'	1:CA:1318:A:O4'	1.86	0.74
2:CC:116:ALA:O	2:CC:119:ILE:HG22	1.87	0.74
16:CS:32:THR:HG22	16:CS:33:TRP:H	1.52	0.74
47:DF:109:ARG:HG3	47:DF:137:PHE:HB3	1.69	0.74
48:DG:132:LEU:HD23	48:DG:132:LEU:H	1.50	0.74
52:DW:24:ARG:HA	52:DW:66:VAL:H	1.52	0.74
23:BB:1579:A:H2'	23:BB:1580:A:H8	1.53	0.74
41:BJ:72:LYS:HA	41:BJ:72:LYS:HE2	1.68	0.74
51:BZ:40:VAL:HG21	51:BZ:43:GLU:HB3	1.67	0.74
16:CS:10:ILE:HG12	16:CS:40:PHE:HZ	1.49	0.74
25:DC:191:LEU:HD23	25:DC:192:GLY:H	1.51	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1011:C:H2'	1:AA:1012:A:H8	1.52	0.74
8:AI:123:ARG:HB2	8:AI:123:ARG:HH11	1.50	0.74
23:BB:30:G:H2'	23:BB:31:C:C6	2.22	0.74
25:BC:4:LYS:HD2	25:BC:5:CYS:H	1.52	0.74
24:DI:105:LEU:HD13	24:DI:129:GLU:HG2	1.67	0.74
23:DB:873:C:H4'	38:DM:64:TRP:HE1	1.51	0.74
46:BU:51:LEU:H	46:BU:53:GLN:HE22	1.35	0.74
23:BB:300:A:H3'	46:BU:81:ARG:HH12	1.51	0.74
1:CA:373:A:H1'	1:CA:481:G:N3	2.02	0.74
5:CF:62:MET:HG3	5:CF:64:VAL:HG23	1.69	0.74
14:CQ:13:SER:HB3	14:CQ:21:VAL:HB	1.68	0.74
26:DD:55:LYS:HZ2	26:DD:60:VAL:HG13	1.53	0.74
24:DI:72:THR:HG21	24:DI:112:LYS:HA	1.69	0.74
23:BB:873:C:H4'	38:BM:64:TRP:HE1	1.53	0.74
23:BB:1205:A:N6	29:BE:165:HIS:HB2	1.95	0.74
29:BE:2:GLU:HA	29:BE:13:THR:HA	1.68	0.74
29:BE:3:LEU:HB2	29:BE:12:LEU:HD12	1.70	0.74
40:BH:31:VAL:O	40:BH:33:GLN:N	2.21	0.74
1:CA:406:G:H21	3:CD:115:GLN:NE2	1.86	0.74
13:CP:28:ARG:HD2	13:CP:29:ASN:N	2.00	0.74
40:DH:14:SER:HB2	40:DH:17:ASP:HB2	1.69	0.74
41:DJ:72:LYS:HE2	41:DJ:72:LYS:HA	1.69	0.74
1:AA:1367:C:H5''	8:AI:115:VAL:HG23	1.68	0.74
47:BF:111:ARG:O	47:BF:112:ASP:HB2	1.87	0.74
24:BI:27:LEU:HD12	24:BI:32:VAL:HG11	1.69	0.74
24:BI:77:VAL:HA	24:BI:80:LYS:HE2	1.70	0.74
4:CE:152:VAL:HG21	7:CH:98:LEU:HB3	1.70	0.74
5:CF:42:TRP:HB2	5:CF:59:TYR:HB2	1.68	0.74
10:CK:19:VAL:HG23	10:CK:34:THR:HG23	1.70	0.74
23:DB:1082:U:N3	23:DB:1086:A:C2	2.55	0.74
23:DB:1796:U:H2'	23:DB:1797:G:H8	1.52	0.74
23:DB:2:G:H2'	23:DB:3:U:C6	2.22	0.74
29:DE:103:GLY:HA2	29:DE:106:LYS:HB3	1.70	0.74
43:DO:66:GLY:HA3	43:DO:102:ARG:NH2	2.03	0.74
1:AA:1412:C:H2'	1:AA:1413:A:C8	2.23	0.74
1:AA:243:A:H4'	1:AA:244:U:H5'	1.69	0.74
1:AA:662:U:H2'	1:AA:663:A:C8	2.21	0.74
9:AJ:52:LEU:H	9:AJ:52:LEU:HD12	1.53	0.74
22:BA:104:A:H2'	22:BA:105:G:O4'	1.88	0.74
25:BC:77:VAL:HG23	25:BC:112:GLY:N	2.03	0.74
19:CU:3:ILE:HB	19:CU:19:LYS:HE3	1.67	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1579:A:H2'	23:DB:1580:A:H8	1.50	0.74
23:DB:593:U:H2'	23:DB:594:U:C6	2.23	0.74
40:DH:48:GLU:HA	40:DH:51:ARG:HE	1.52	0.74
50:DT:69:ARG:HE	50:DT:69:ARG:HA	1.53	0.74
1:AA:1406:U:H2'	1:AA:1407:C:H5'	1.69	0.74
3:AD:84:ASN:ND2	4:AE:101:GLY:HA3	2.02	0.74
8:AI:18:VAL:HG11	8:AI:82:ILE:HA	1.70	0.74
26:BD:151:THR:HB	26:BD:152:PRO:HD3	1.68	0.74
44:BQ:16:ILE:HA	44:BQ:19:GLN:HB2	1.68	0.74
2:CC:149:LYS:HB3	2:CC:200:TRP:HB2	1.69	0.74
23:DB:1099:G:O4'	24:DI:3:LYS:C	2.25	0.74
23:DB:30:G:H2'	23:DB:31:C:C6	2.22	0.74
23:DB:590:A:H2'	23:DB:591:U:C6	2.23	0.74
26:DD:186:LEU:HD11	28:DP:3:ILE:HG13	1.69	0.74
1:AA:310:G:H5''	13:AP:31:ARG:HB2	1.69	0.74
37:BL:93:ASN:ND2	37:BL:94:THR:H	1.86	0.74
1:CA:406:G:H21	3:CD:115:GLN:HE22	1.36	0.74
18:CB:95:TRP:HH2	18:CB:100:LEU:HB2	1.52	0.74
12:CM:15:VAL:HA	12:CM:29:SER:OG	1.88	0.74
23:DB:1141:U:H5''	41:DJ:27:ARG:HH21	1.53	0.74
25:DC:140:VAL:HG12	25:DC:141:HIS:H	1.52	0.74
3:AD:26:ALA:HA	3:AD:30:LYS:HE3	1.70	0.74
10:AK:91:GLY:O	10:AK:95:THR:HG22	1.88	0.74
11:AL:14:LYS:HG2	11:AL:16:ALA:H	1.53	0.74
16:AS:44:ILE:HG13	16:AS:62:THR:HA	1.69	0.74
23:BB:2848:G:H1	23:BB:2867:G:N2	1.85	0.74
1:CA:1241:G:H2'	1:CA:1242:G:H8	1.53	0.74
1:AA:337:G:H2'	1:AA:338:A:C8	2.22	0.73
23:BB:2147:A:H5'	23:BB:2148:G:O4'	1.88	0.73
26:BD:9:VAL:HA	26:BD:197:THR:HG23	1.70	0.73
40:BH:4:ILE:HB	40:BH:37:VAL:HG12	1.69	0.73
37:DL:80:SER:HA	37:DL:115:GLU:HB2	1.68	0.73
6:CG:145:GLU:CA	6:CG:148:LYS:HB2	2.18	0.73
23:DB:1729:U:H3'	23:DB:1730:C:O4'	1.87	0.73
23:DB:2386:A:H4'	52:DW:54:ARG:O	1.88	0.73
23:DB:354:A:H2'	23:DB:355:U:C6	2.23	0.73
23:DB:580:U:H2'	23:DB:581:C:C6	2.23	0.73
26:DD:116:LYS:HG3	26:DD:165:MET:SD	2.28	0.73
47:DF:111:ARG:O	47:DF:112:ASP:HB2	1.88	0.73
47:DF:147:ARG:HG2	47:DF:148:VAL:HG22	1.69	0.73
28:DP:103:THR:HG22	28:DP:104:GLY:H	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:94:ARG:HH11	8:AI:94:ARG:HB3	1.53	0.73
12:AM:21:ILE:HG22	12:AM:23:GLY:H	1.53	0.73
33:B1:33:LEU:HB3	33:B1:51:ALA:CB	2.18	0.73
7:CH:94:VAL:HG23	7:CH:101:ALA:HB2	1.70	0.73
23:DB:101:A:O2'	23:DB:102:U:H5''	1.88	0.73
25:DC:77:VAL:HG23	25:DC:112:GLY:N	2.02	0.73
29:DE:148:ILE:HG13	29:DE:167:VAL:HG23	1.70	0.73
47:DF:41:GLU:HB3	47:DF:44:ALA:HB3	1.71	0.73
47:DF:8:LYS:HA	47:DF:12:VAL:HG21	1.67	0.73
45:DS:24:ILE:HG23	45:DS:32:ALA:HB1	1.70	0.73
2:AC:63:ILE:HD12	2:AC:98:ALA:HB2	1.70	0.73
23:BB:184:C:H2'	23:BB:185:G:H8	1.52	0.73
23:BB:2512:C:H2'	23:BB:2513:A:O4'	1.88	0.73
47:BF:11:VAL:HG12	47:BF:12:VAL:H	1.52	0.73
47:BF:146:ASP:HB3	47:BF:147:ARG:HH12	1.51	0.73
2:CC:8:GLY:HA2	2:CC:11:LEU:HG	1.70	0.73
23:DB:2186:G:H2'	23:DB:2187:U:O4'	1.87	0.73
48:DG:96:ALA:HB3	48:DG:103:ASN:HB3	1.68	0.73
2:AC:120:THR:HG21	2:AC:186:SER:HB3	1.71	0.73
23:BB:1796:U:H2'	23:BB:1797:G:H8	1.53	0.73
47:BF:111:ARG:NH2	47:BF:113:PHE:HB2	2.03	0.73
23:BB:2305:U:H5''	47:BF:130:GLY:HA3	1.70	0.73
41:BJ:55:ILE:HG22	41:BJ:123:LYS:HB2	1.70	0.73
23:DB:2153:C:H2'	23:DB:2154:A:H8	1.52	0.73
38:DM:40:ARG:HD3	38:DM:93:VAL:HG21	1.70	0.73
9:AJ:36:VAL:HG12	9:AJ:38:GLY:H	1.53	0.73
10:AK:22:ILE:HB	10:AK:85:VAL:HG22	1.69	0.73
1:CA:859:G:H2'	1:CA:860:A:C8	2.24	0.73
1:CA:977:A:H2'	1:CA:978:A:H5''	1.69	0.73
18:CB:102:ASN:ND2	18:CB:105:THR:HB	2.02	0.73
21:CN:27:LYS:HG3	21:CN:28:ALA:H	1.54	0.73
34:D3:49:VAL:HG11	34:D3:54:LEU:HD13	1.71	0.73
23:DB:2872:A:O2'	23:DB:2873:A:H5'	1.88	0.73
1:AA:337:G:H2'	1:AA:338:A:H8	1.52	0.73
1:AA:699:C:H2'	1:AA:700:G:H5''	1.71	0.73
1:AA:9:G:H5'	4:AE:107:GLY:HA3	1.70	0.73
23:BB:593:U:H2'	23:BB:594:U:C6	2.24	0.73
40:BH:14:SER:HB2	40:BH:17:ASP:HB2	1.71	0.73
2:CC:113:LYS:HD2	2:CC:184:ASN:ND2	2.03	0.73
13:CP:67:ILE:HD11	13:CP:71:VAL:HG22	1.71	0.73
14:CQ:20:ILE:HD13	14:CQ:47:ASP:HB3	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CS:15:LEU:HA	16:CS:18:VAL:HG12	1.70	0.73
23:DB:2776:A:H4'	23:DB:2777:G:H5''	1.71	0.73
27:DK:19:VAL:HG13	27:DK:43:ILE:HA	1.69	0.73
35:DV:77:VAL:HA	35:DV:89:ILE:HG22	1.69	0.73
1:AA:462:G:N3	1:AA:462:G:H2'	2.03	0.73
4:AE:73:VAL:HG21	4:AE:143:LEU:HB3	1.70	0.73
8:AI:61:ASP:O	8:AI:62:LEU:HD13	1.89	0.73
12:AM:21:ILE:HB	12:AM:24:VAL:HG22	1.71	0.73
29:BE:158:PHE:HA	29:BE:169:VAL:HG21	1.71	0.73
40:BH:58:LEU:O	40:BH:62:LEU:HG	1.87	0.73
49:BR:28:ALA:O	49:BR:63:VAL:HG21	1.89	0.73
1:CA:662:U:H2'	1:CA:663:A:C8	2.23	0.73
18:CB:122:ASP:OD1	18:CB:124:THR:HG22	1.89	0.73
2:CC:5:HIS:CD2	2:CC:8:GLY:H	2.06	0.73
21:CN:65:GLN:HB2	21:CN:82:LYS:HE2	1.70	0.73
9:CJ:66:GLU:N	21:CN:98:ALA:HB2	2.04	0.73
40:DH:86:ASP:HB2	40:DH:89:LYS:HB2	1.69	0.73
42:DN:97:ILE:HD12	42:DN:98:LEU:N	2.03	0.73
35:DV:38:LEU:HG	35:DV:40:ILE:HD12	1.69	0.73
1:AA:1317:C:OP1	21:AN:56:PRO:HD2	1.89	0.73
23:BB:191:A:H2'	23:BB:192:C:C6	2.24	0.73
6:CG:68:VAL:HG11	6:CG:103:ILE:HG13	1.70	0.73
9:CJ:57:VAL:HG22	9:CJ:58:ASN:H	1.52	0.73
11:CL:35:ARG:HH21	11:CL:36:VAL:HG22	1.53	0.73
37:DL:57:LEU:HD12	37:DL:60:ARG:HH22	1.53	0.73
1:AA:961:U:OP2	1:AA:1223:C:H1'	1.89	0.73
2:AC:147:GLY:HA3	2:AC:202:PHE:HB3	1.71	0.73
5:AF:19:PRO:HA	5:AF:22:ILE:HD12	1.70	0.73
7:AH:103:VAL:HG22	7:AH:124:ILE:HA	1.70	0.73
7:AH:11:THR:HG22	7:AH:14:ARG:HH12	1.52	0.73
13:AP:28:ARG:CD	13:AP:29:ASN:H	2.02	0.73
23:BB:1124:G:H1'	32:B4:38:GLY:OXT	1.87	0.73
50:BT:48:GLN:HE22	50:BT:55:VAL:HB	1.53	0.73
1:CA:859:G:H2'	1:CA:860:A:H8	1.54	0.73
25:DC:183:VAL:HG13	25:DC:185:ALA:H	1.54	0.73
29:DE:105:LEU:HA	29:DE:108:ILE:HG22	1.71	0.73
26:DD:118:PHE:CE2	42:DN:1:MET:HB3	2.24	0.73
5:AF:62:MET:HG3	5:AF:64:VAL:HG23	1.70	0.72
9:AJ:8:ILE:HD13	9:AJ:75:ASP:HA	1.71	0.72
21:AN:50:LEU:H	21:AN:51:PRO:CD	2.02	0.72
31:D0:29:VAL:HG23	31:D0:35:GLU:H	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:91:THR:HG23	26:DD:92:VAL:H	1.54	0.72
29:DE:158:PHE:HA	29:DE:169:VAL:HG21	1.69	0.72
44:DQ:63:ARG:HH21	44:DQ:64:ILE:HD11	1.54	0.72
46:DU:82:VAL:H	46:DU:96:LYS:HZ2	1.36	0.72
1:AA:1499:A:H5'	1:AA:1499:A:H8	1.52	0.72
18:AB:202:ASN:HD22	18:AB:204:ASP:N	1.82	0.72
23:BB:2267:A:C8	23:BB:2267:A:H3'	2.23	0.72
23:BB:877:A:H2'	23:BB:899:A:N1	2.04	0.72
1:CA:1290:G:H2'	1:CA:1291:U:C6	2.22	0.72
1:CA:390:U:H2'	1:CA:391:G:H8	1.54	0.72
1:CA:674:G:H2'	1:CA:675:A:H8	1.52	0.72
23:DB:2346:A:H3'	23:DB:2347:C:H5''	1.70	0.72
23:DB:2800:A:H2'	23:DB:2801:G:O4'	1.89	0.72
23:DB:477:A:H2'	23:DB:478:A:C8	2.24	0.72
23:DB:532:A:N3	23:DB:532:A:H2'	2.04	0.72
4:AE:152:VAL:HG21	7:AH:98:LEU:HB3	1.69	0.72
22:BA:49:C:H2'	22:BA:50:A:H8	1.55	0.72
23:BB:1324:G:H1'	23:BB:1616:A:N6	2.04	0.72
23:BB:1729:U:H3'	23:BB:1730:C:O4'	1.89	0.72
23:BB:858:G:H21	23:BB:2268:A:H3'	1.53	0.72
47:BF:128:SER:HB3	47:BF:154:THR:HG23	1.71	0.72
44:BQ:60:TRP:O	44:BQ:64:ILE:HG12	1.89	0.72
44:BQ:90:ASP:O	44:BQ:94:LEU:HB2	1.89	0.72
44:BQ:93:ILE:HG23	44:BQ:94:LEU:HD22	1.70	0.72
35:BV:77:VAL:HA	35:BV:89:ILE:HG22	1.69	0.72
21:CN:25:GLU:HB2	21:CN:29:ILE:HD11	1.70	0.72
23:DB:698:C:H4'	23:DB:734:A:H61	1.53	0.72
41:DJ:6:ALA:HB3	41:DJ:45:THR:HG21	1.71	0.72
28:DP:50:ARG:HD3	28:DP:56:SER:HB3	1.71	0.72
44:DQ:60:TRP:O	44:DQ:64:ILE:HG12	1.89	0.72
31:B0:29:VAL:HG23	31:B0:35:GLU:H	1.53	0.72
23:BB:135:U:H2'	23:BB:136:G:C8	2.24	0.72
23:BB:1843:C:H5''	25:BC:250:GLN:NE2	2.04	0.72
23:BB:2064:C:H2'	23:BB:2065:C:C6	2.24	0.72
23:BB:544:C:H2'	23:BB:545:U:O2	1.89	0.72
48:BG:84:LYS:HB3	48:BG:132:LEU:O	1.90	0.72
52:BW:24:ARG:HA	52:BW:66:VAL:H	1.52	0.72
22:DA:104:A:H2'	22:DA:105:G:O4'	1.89	0.72
24:DI:20:SER:HB3	24:DI:21:PRO:HD3	1.69	0.72
1:AA:238:A:H2'	1:AA:239:U:H5''	1.70	0.72
1:AA:269:C:H2'	1:AA:270:A:H8	1.52	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:62:SER:HA	2:AC:97:PRO:HG2	1.71	0.72
5:AF:1:MET:SD	5:AF:67:PRO:HD3	2.29	0.72
8:AI:41:GLU:O	8:AI:44:ARG:HG2	1.89	0.72
23:BB:2820:A:OP1	42:BN:4:ARG:HA	1.90	0.72
23:BB:2872:A:O2'	23:BB:2873:A:H5'	1.88	0.72
23:BB:477:A:H2'	23:BB:478:A:C8	2.24	0.72
25:BC:140:VAL:HG12	25:BC:141:HIS:H	1.53	0.72
1:CA:203:G:H1'	1:CA:465:A:C6	2.25	0.72
1:CA:946:A:H2'	1:CA:947:G:H8	1.53	0.72
12:CM:3:ILE:HA	12:CM:56:ARG:HG3	1.72	0.72
22:DA:2:G:H2'	22:DA:3:C:H6	1.53	0.72
23:DB:1412:U:H2'	23:DB:1413:A:C8	2.24	0.72
23:DB:773:U:H5'	23:DB:774:G:OP2	1.89	0.72
25:DC:16:VAL:H	25:DC:203:VAL:HG11	1.53	0.72
44:DQ:16:ILE:HA	44:DQ:19:GLN:HB2	1.70	0.72
1:AA:1342:C:H2'	1:AA:1343:G:H8	1.52	0.72
1:AA:176:C:H2'	1:AA:177:G:N3	2.04	0.72
7:AH:49:LYS:HB3	7:AH:59:GLU:HB2	1.71	0.72
12:AM:90:HIS:HA	12:AM:108:ARG:NH2	2.05	0.72
33:B1:34:GLU:HG2	33:B1:49:LYS:HG3	1.70	0.72
47:BF:41:GLU:HB3	47:BF:44:ALA:HB3	1.71	0.72
42:BN:97:ILE:HD12	42:BN:98:LEU:N	2.04	0.72
1:CA:699:C:H2'	1:CA:700:G:H5''	1.72	0.72
18:CB:63:LYS:HA	18:CB:224:ARG:NH1	2.05	0.72
23:DB:1080:A:H4'	24:DI:126:ARG:HD3	1.71	0.72
1:AA:1048:G:N2	1:AA:1214:C:H5	1.86	0.72
23:BB:2346:A:H3'	23:BB:2347:C:H5''	1.71	0.72
23:BB:571:U:H3'	49:BR:80:ARG:NH1	2.04	0.72
1:CA:1101:A:H62	18:CB:173:LYS:HE3	1.54	0.72
1:CA:977:A:C2	1:CA:1223:C:H2'	2.25	0.72
5:CF:6:ILE:HD11	5:CF:8:PHE:HD2	1.54	0.72
17:CT:2:ASN:ND2	17:CT:3:ILE:HG13	2.05	0.72
23:DB:1174:U:O2'	23:DB:1175:A:H3'	1.89	0.72
23:DB:117:G:H5'	23:DB:126:A:H8	1.54	0.72
23:DB:1937:A:N7	23:DB:1939:U:H2'	2.05	0.72
23:DB:2092:U:H4'	23:DB:2093:G:O5'	1.88	0.72
25:DC:132:ARG:HH11	25:DC:186:ASP:HB2	1.55	0.72
26:DD:5:VAL:N	26:DD:32:ASN:HD21	1.86	0.72
29:DE:105:LEU:HD11	29:DE:177:PRO:HG3	1.72	0.72
35:DV:72:VAL:HG21	35:DV:91:PHE:HB3	1.69	0.72
1:AA:1011:C:H2'	1:AA:1012:A:C8	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:406:G:H21	3:AD:115:GLN:NE2	1.88	0.72
2:AC:87:ARG:HH21	2:AC:88:LYS:HA	1.55	0.72
21:AN:64:ARG:HB2	21:AN:77:GLY:O	1.89	0.72
13:AP:67:ILE:HD11	13:AP:71:VAL:HG22	1.70	0.72
23:BB:1287:A:H3'	23:BB:1288:G:N2	2.05	0.72
23:BB:2502:G:H5'	23:BB:2503:A:H5''	1.72	0.72
23:BB:674:G:O3'	29:BE:60:TRP:HZ2	1.72	0.72
1:AA:1432:G:C5'	28:BP:105:LYS:HG2	2.19	0.72
1:CA:1461:G:H2'	1:CA:1462:C:H6	1.54	0.72
18:CB:80:LYS:HG3	18:CB:81:ASP:H	1.55	0.72
40:DH:31:VAL:O	40:DH:33:GLN:N	2.22	0.72
44:DQ:42:GLY:HA3	49:DR:75:VAL:HG21	1.71	0.72
1:AA:1237:C:H3'	1:AA:1238:A:H5'	1.71	0.72
4:AE:156:ARG:NH1	7:AH:42:GLU:HB3	2.05	0.72
5:AF:6:ILE:HD11	5:AF:8:PHE:HD2	1.55	0.72
23:BB:1412:U:H2'	23:BB:1413:A:H8	1.55	0.72
23:BB:2758:A:H2'	23:BB:2759:G:O4'	1.90	0.72
23:BB:2307:G:O6	47:BF:84:ILE:HD11	1.89	0.72
27:BK:112:PHE:O	27:BK:115:ILE:HG22	1.90	0.72
1:CA:1316:G:N2	1:CA:1318:A:H3'	2.05	0.72
1:CA:649:A:H2'	1:CA:650:G:O4'	1.89	0.72
18:CB:31:PHE:HB2	18:CB:41:ASN:HA	1.72	0.72
19:CU:24:LYS:HZ3	19:CU:25:ALA:H	1.37	0.72
47:DF:11:VAL:HG12	47:DF:12:VAL:H	1.53	0.72
42:DN:97:ILE:HD12	42:DN:98:LEU:H	1.55	0.72
18:AB:120:SER:HA	18:AB:125:PHE:HB3	1.72	0.72
23:BB:532:A:N3	23:BB:532:A:H2'	2.05	0.72
25:BC:191:LEU:HD23	25:BC:192:GLY:H	1.54	0.72
29:BE:103:GLY:HA2	29:BE:106:LYS:HB3	1.72	0.72
48:BG:122:ALA:HA	48:BG:132:LEU:HA	1.72	0.72
40:BH:121:VAL:HB	40:BH:128:HIS:NE2	2.05	0.72
18:CB:59:ILE:HG22	18:CB:62:ARG:HE	1.53	0.72
12:CM:90:HIS:HA	12:CM:108:ARG:NH2	2.04	0.72
23:DB:2074:U:H1'	23:DB:2598:A:N3	2.05	0.72
47:DF:146:ASP:HB3	47:DF:147:ARG:HH12	1.53	0.72
43:DO:27:VAL:HG21	43:DO:40:ILE:HD12	1.71	0.72
50:DT:48:GLN:HE22	50:DT:55:VAL:HB	1.54	0.72
5:AF:42:TRP:HB2	5:AF:59:TYR:HB2	1.71	0.71
23:BB:2800:A:H2'	23:BB:2801:G:O4'	1.89	0.71
35:BV:44:HIS:HE1	35:BV:86:LEU:H	1.38	0.71
1:CA:1032:G:H2'	1:CA:1033:G:O4'	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CK:91:GLY:O	10:CK:95:THR:HG22	1.90	0.71
23:DB:1060:U:C4	23:DB:1088:A:N6	2.58	0.71
23:DB:1871:A:H2'	23:DB:1872:A:C8	2.25	0.71
23:DB:899:A:H3'	23:DB:900:A:C8	2.25	0.71
40:DH:65:ALA:HB1	40:DH:138:VAL:HG21	1.72	0.71
27:DK:19:VAL:HG12	27:DK:41:ILE:HG12	1.71	0.71
1:AA:1053:G:N7	1:AA:1199:U:H3'	2.04	0.71
1:AA:649:A:H2'	1:AA:650:G:O4'	1.90	0.71
11:AL:35:ARG:HH21	11:AL:36:VAL:HG22	1.55	0.71
20:AO:11:ILE:HG12	20:AO:30:ALA:HB1	1.70	0.71
16:AS:31:ARG:HG3	16:AS:56:HIS:NE2	2.05	0.71
33:B1:47:ILE:H	33:B1:47:ILE:HD12	1.54	0.71
23:BB:1412:U:H2'	23:BB:1413:A:C8	2.25	0.71
23:BB:1683:U:H2'	23:BB:1684:G:C8	2.25	0.71
47:BF:109:ARG:HG3	47:BF:137:PHE:HB3	1.71	0.71
40:BH:147:VAL:HG12	40:BH:148:ALA:H	1.54	0.71
42:BN:101:GLY:HA2	42:BN:109:PRO:HA	1.71	0.71
37:DL:93:ASN:ND2	37:DL:94:THR:H	1.88	0.71
49:DR:28:ALA:O	49:DR:63:VAL:HG21	1.89	0.71
17:AT:2:ASN:ND2	17:AT:3:ILE:HG13	2.04	0.71
19:AU:3:ILE:HB	19:AU:19:LYS:HE3	1.71	0.71
23:BB:1889:A:H2'	23:BB:1890:A:C8	2.25	0.71
23:BB:698:C:H4'	23:BB:734:A:H61	1.54	0.71
24:BI:20:SER:HB3	24:BI:21:PRO:HD3	1.71	0.71
27:BK:71:ARG:CZ	27:BK:72:PRO:HD3	2.20	0.71
52:BW:39:GLN:HG3	52:BW:42:THR:CB	2.20	0.71
1:CA:988:G:H21	1:CA:1015:G:N2	1.87	0.71
23:DB:1098:A:H2'	24:DI:4:VAL:N	2.04	0.71
23:DB:2291:U:H2'	23:DB:2292:U:C6	2.25	0.71
23:DB:661:A:H1'	37:DL:12:SER:O	1.90	0.71
1:AA:190:A:H2'	1:AA:191:G:O4'	1.91	0.71
1:AA:394:G:H2'	1:AA:395:C:H6	1.55	0.71
1:AA:920:U:H2'	1:AA:921:U:C6	2.26	0.71
23:BB:1911:U:H2'	23:BB:1918:A:N1	2.05	0.71
29:BE:75:SER:O	29:BE:78:TRP:HB2	1.90	0.71
43:BO:66:GLY:HA3	43:BO:102:ARG:NH2	2.04	0.71
44:BQ:63:ARG:HH21	44:BQ:64:ILE:HD11	1.56	0.71
1:CA:462:G:H2'	1:CA:462:G:N3	2.04	0.71
3:CD:29:THR:HG22	3:CD:30:LYS:H	1.55	0.71
4:CE:73:VAL:HG21	4:CE:143:LEU:HB3	1.71	0.71
23:DB:184:C:H2'	23:DB:185:G:H8	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2153:C:H2'	23:DB:2154:A:C8	2.25	0.71
30:DY:3:THR:HB	30:DY:36:GLU:HG3	1.72	0.71
1:AA:1342:C:H5'	8:AI:127:SER:HA	1.73	0.71
23:BB:322:A:OP2	29:BE:163:ASN:HB2	1.90	0.71
1:CA:449:G:H2'	1:CA:450:G:C8	2.25	0.71
15:CR:26:ALA:HA	15:CR:29:LYS:HE2	1.73	0.71
22:DA:11:C:H3'	22:DA:12:C:H5''	1.73	0.71
22:DA:54:G:H21	47:DF:25:MET:HG2	1.55	0.71
23:DB:1098:A:H2'	24:DI:4:VAL:CA	2.20	0.71
23:DB:704:G:H2'	23:DB:726:G:N2	2.04	0.71
27:DK:118:LEU:C	27:DK:120:PRO:HD2	2.10	0.71
51:DZ:56:MET:HA	51:DZ:59:ILE:HG12	1.73	0.71
9:AJ:40:ILE:HD13	9:AJ:73:LEU:HD12	1.72	0.71
23:BB:1141:U:H5''	41:BJ:27:ARG:HH21	1.53	0.71
29:BE:105:LEU:HA	29:BE:108:ILE:HG22	1.73	0.71
24:BI:106:GLN:O	24:BI:110:GLN:HG3	1.90	0.71
50:BT:60:THR:HB	50:BT:81:LYS:HD2	1.73	0.71
52:BW:35:ILE:HA	52:BW:57:THR:HG23	1.71	0.71
8:CI:24:ASN:HA	8:CI:58:GLU:HA	1.73	0.71
21:CN:5:MET:HA	21:CN:8:ARG:HG3	1.72	0.71
33:D1:37:LYS:H	33:D1:48:TYR:HD2	1.36	0.71
23:DB:1412:U:H2'	23:DB:1413:A:H8	1.56	0.71
23:DB:2548:U:H1'	27:DK:23:LYS:NZ	2.04	0.71
1:AA:390:U:H2'	1:AA:391:G:H8	1.55	0.71
3:AD:12:ARG:HH11	3:AD:12:ARG:HB3	1.55	0.71
12:AM:2:ARG:O	12:AM:7:ASN:HB2	1.89	0.71
22:BA:43:C:H1'	47:BF:91:ARG:HD2	1.71	0.71
23:BB:950:G:H2'	23:BB:951:C:C6	2.26	0.71
48:BG:9:VAL:HA	48:BG:48:THR:HB	1.72	0.71
24:BI:122:GLU:O	24:BI:126:ARG:HG3	1.90	0.71
28:BP:103:THR:HG22	28:BP:104:GLY:H	1.54	0.71
1:CA:269:C:H2'	1:CA:270:A:H8	1.53	0.71
9:CJ:6:ILE:HD12	9:CJ:76:ILE:HD11	1.73	0.71
33:D1:47:ILE:H	33:D1:47:ILE:HD12	1.54	0.71
23:DB:1459:G:H2'	23:DB:1461:C:C5	2.25	0.71
23:DB:1889:A:H2'	23:DB:1890:A:C8	2.26	0.71
1:AA:552:U:H5'	11:AL:82:ARG:HH11	1.55	0.71
4:AE:155:LYS:HG2	7:AH:70:VAL:HG13	1.73	0.71
16:AS:52:ASN:CG	16:AS:53:GLY:H	1.94	0.71
23:BB:1812:U:H2'	23:BB:1813:G:H8	1.55	0.71
23:BB:2008:C:H2'	23:BB:2009:A:H8	1.56	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:773:U:H5'	23:BB:774:G:OP2	1.91	0.71
23:BB:1843:C:H5''	25:BC:250:GLN:HE21	1.55	0.71
48:BG:148:ARG:HD3	48:BG:152:ARG:HH11	1.55	0.71
42:BN:38:LEU:HB3	42:BN:39:PRO:HD3	1.73	0.71
8:CI:66:VAL:CG2	8:CI:74:GLN:HG3	2.20	0.71
21:CN:68:ARG:NH1	21:CN:70:HIS:HB2	2.04	0.71
47:DF:111:ARG:NH2	47:DF:113:PHE:HB2	2.06	0.71
47:DF:128:SER:HB3	47:DF:154:THR:HG23	1.73	0.71
43:DO:83:LEU:HD11	43:DO:114:GLY:HA3	1.73	0.71
8:AI:11:ARG:H	8:AI:14:SER:HB3	1.56	0.71
33:B1:9:LYS:H	33:B1:9:LYS:HD3	1.56	0.71
23:BB:1060:U:C4	23:BB:1088:A:N6	2.58	0.71
49:BR:72:VAL:HG23	49:BR:89:HIS:HB3	1.72	0.71
1:CA:1144:G:N2	1:CA:1146:A:H62	1.89	0.71
1:CA:1521:C:H2'	1:CA:1522:U:C6	2.25	0.71
4:CE:133:ILE:HD12	4:CE:133:ILE:H	1.56	0.71
5:CF:19:PRO:HA	5:CF:22:ILE:HD12	1.71	0.71
20:CO:70:LEU:HD11	20:CO:77:ARG:HB2	1.72	0.71
33:D1:5:ARG:HH11	33:D1:25:ASN:HB2	1.56	0.71
23:DB:659:G:H21	29:DE:30:GLN:HE22	1.36	0.71
27:DK:71:ARG:CZ	27:DK:72:PRO:HD3	2.20	0.71
44:DQ:93:ILE:HG23	44:DQ:94:LEU:HD22	1.71	0.71
1:AA:1089:G:H2'	1:AA:1090:U:H5'	1.73	0.71
23:BB:1141:U:H4'	23:BB:1142:A:O4'	1.89	0.71
23:BB:28:A:H61	23:BB:512:G:H1'	1.56	0.71
25:BC:123:ILE:HD12	25:BC:191:LEU:HD11	1.71	0.71
41:BJ:3:THR:HG21	44:BQ:60:TRP:HE1	1.55	0.71
1:CA:552:U:H5'	11:CL:82:ARG:HH11	1.56	0.71
1:CA:946:A:H2'	1:CA:947:G:C8	2.26	0.71
13:CP:28:ARG:CD	13:CP:29:ASN:H	2.01	0.71
32:D4:3:VAL:HG23	32:D4:4:ARG:H	1.56	0.71
23:DB:1324:G:H1'	23:DB:1616:A:N6	2.04	0.71
47:DF:66:ILE:HD12	47:DF:85:GLY:O	1.91	0.71
23:DB:1287:A:OP1	42:DN:104:ALA:HB3	1.91	0.71
15:AR:72:ARG:HH11	15:AR:72:ARG:N	1.86	0.70
40:BH:116:ARG:NH2	40:BH:139:PHE:HB3	2.06	0.70
1:CA:1347:G:N2	1:CA:1373:G:H2'	2.06	0.70
1:CA:394:G:H2'	1:CA:395:C:H6	1.55	0.70
1:CA:87:C:H2'	1:CA:88:U:H5'	1.73	0.70
2:CC:6:PRO:HB3	2:CC:174:LEU:HD23	1.73	0.70
12:CM:21:ILE:HG23	12:CM:65:GLU:OE2	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CS:10:ILE:HG23	16:CS:37:SER:HB3	1.73	0.70
23:DB:1548:A:H2'	23:DB:1549:A:C8	2.25	0.70
42:DN:101:GLY:HA2	42:DN:110:MET:N	2.05	0.70
52:DW:39:GLN:HG3	52:DW:42:THR:CB	2.21	0.70
1:AA:950:U:H2'	1:AA:951:G:H8	1.57	0.70
2:AC:63:ILE:HD11	2:AC:96:VAL:HG23	1.73	0.70
23:BB:2886:A:H62	31:B0:39:ARG:CZ	2.03	0.70
26:BD:111:GLY:H	26:BD:194:PRO:HG2	1.54	0.70
42:BN:101:GLY:HA2	42:BN:110:MET:N	2.05	0.70
1:CA:437:U:H2'	1:CA:438:U:O4'	1.91	0.70
18:CB:56:LEU:O	18:CB:60:ALA:HB2	1.91	0.70
23:DB:1000:A:H2'	23:DB:1001:A:C8	2.25	0.70
23:DB:1138:G:H2'	23:DB:1139:G:O4'	1.92	0.70
23:DB:832:U:H2'	23:DB:833:A:C8	2.26	0.70
24:DI:85:ILE:HD13	24:DI:137:LEU:HD21	1.73	0.70
11:AL:86:VAL:HG23	11:AL:89:LEU:HB2	1.72	0.70
23:BB:2458:G:H8	23:BB:2459:A:H62	1.40	0.70
23:BB:2824:C:H3'	23:BB:2825:G:H21	1.56	0.70
23:BB:287:G:H2'	23:BB:288:U:H6	1.55	0.70
37:BL:90:VAL:HB	37:BL:122:VAL:HG12	1.72	0.70
4:CE:155:LYS:HG2	7:CH:70:VAL:HG13	1.73	0.70
8:CI:105:ARG:HD3	8:CI:106:ASP:N	2.06	0.70
15:CR:40:PRO:HD2	15:CR:43:ILE:HD11	1.71	0.70
23:DB:264:C:H2'	23:DB:265:A:H5''	1.71	0.70
25:DC:68:ARG:HH21	25:DC:190:THR:HG22	1.56	0.70
50:DT:66:LYS:O	50:DT:77:ARG:HB2	1.91	0.70
46:DU:51:LEU:H	46:DU:53:GLN:HE22	1.38	0.70
1:AA:301:G:H2'	1:AA:302:G:H8	1.54	0.70
1:AA:41:G:H2'	1:AA:42:G:H8	1.55	0.70
1:AA:715:A:H2'	1:AA:716:A:C8	2.27	0.70
1:AA:967:C:H3'	1:AA:968:A:H5'	1.71	0.70
23:BB:1000:A:H2'	23:BB:1001:A:C8	2.25	0.70
23:BB:2748:A:H1'	48:BG:66:THR:HB	1.73	0.70
48:BG:84:LYS:HG3	48:BG:131:VAL:HB	1.73	0.70
27:BK:10:VAL:HG13	27:BK:86:LEU:HD21	1.73	0.70
1:CA:1360:A:H2'	1:CA:1361:G:O4'	1.90	0.70
1:CA:940:C:H2'	1:CA:941:G:C8	2.26	0.70
23:DB:1553:A:O2'	23:DB:1554:U:H2'	1.91	0.70
29:DE:63:LYS:HE3	29:DE:63:LYS:HA	1.73	0.70
48:DG:122:ALA:HA	48:DG:132:LEU:HA	1.71	0.70
27:DK:88:ASN:HD22	27:DK:89:ASN:N	1.88	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1057:G:H2'	1:AA:1058:G:O4'	1.91	0.70
1:AA:1369:C:H2'	1:AA:1370:G:C8	2.26	0.70
23:BB:1287:A:OP1	42:BN:104:ALA:HB3	1.91	0.70
23:BB:2776:A:H4'	23:BB:2777:G:H5''	1.73	0.70
27:BK:88:ASN:HD22	27:BK:89:ASN:N	1.89	0.70
38:BM:40:ARG:HD3	38:BM:93:VAL:HG21	1.74	0.70
28:BP:50:ARG:HD3	28:BP:56:SER:HB3	1.73	0.70
50:BT:66:LYS:O	50:BT:77:ARG:HB2	1.91	0.70
35:BV:51:GLN:HB2	35:BV:57:TYR:OH	1.90	0.70
1:CA:238:A:H2'	1:CA:239:U:H5''	1.74	0.70
1:CA:310:G:H5''	13:CP:31:ARG:HB2	1.72	0.70
1:CA:41:G:H2'	1:CA:42:G:H8	1.54	0.70
12:CM:28:ARG:NH2	12:CM:62:PHE:HB2	2.05	0.70
23:DB:1381:G:H2'	23:DB:1382:G:H5'	1.74	0.70
23:DB:1854:A:N6	23:DB:1888:G:H1'	2.06	0.70
26:DD:159:LYS:HZ2	26:DD:160:LYS:N	1.90	0.70
43:DO:67:ASN:HB3	43:DO:70:ALA:HB2	1.73	0.70
51:DZ:64:ILE:HD12	51:DZ:64:ILE:H	1.55	0.70
1:AA:449:G:H2'	1:AA:450:G:C8	2.27	0.70
1:AA:674:G:H2'	1:AA:675:A:H8	1.54	0.70
13:AP:57:ILE:O	13:AP:61:VAL:HG23	1.91	0.70
14:AQ:74:LEU:HD22	14:AQ:75:VAL:N	2.07	0.70
15:AR:40:PRO:HD2	15:AR:43:ILE:HD11	1.72	0.70
22:BA:32:U:H4'	22:BA:52:A:N6	2.06	0.70
1:CA:239:U:OP1	1:CA:239:U:H4'	1.92	0.70
5:CF:2:ARG:HG3	5:CF:92:THR:OG1	1.92	0.70
23:DB:2098:U:H2'	23:DB:2099:U:O4'	1.90	0.70
26:DD:16:THR:HB	26:DD:18:ASP:OD1	1.91	0.70
29:DE:75:SER:O	29:DE:78:TRP:HB2	1.92	0.70
47:DF:102:LEU:HA	47:DF:106:ALA:HB2	1.74	0.70
49:DR:68:ARG:NH1	49:DR:90:ARG:HD3	2.06	0.70
1:AA:390:U:H2'	1:AA:391:G:C8	2.27	0.70
10:AK:23:HIS:HB3	10:AK:30:ILE:HG13	1.74	0.70
11:AL:49:ARG:HG2	11:AL:89:LEU:HD21	1.74	0.70
12:AM:102:LYS:HG3	12:AM:103:THR:H	1.55	0.70
23:BB:922:C:H1'	52:BW:22:VAL:HG21	1.72	0.70
25:BC:16:VAL:H	25:BC:203:VAL:HG11	1.56	0.70
28:BP:23:ASP:HA	28:BP:88:ARG:HA	1.73	0.70
1:CA:1250:A:H4'	8:CI:69:GLY:H	1.56	0.70
23:DB:2150:C:H2'	23:DB:2151:U:H6	1.55	0.70
23:DB:2848:G:H1	23:DB:2867:G:N2	1.89	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:950:G:H2'	23:DB:951:C:C6	2.27	0.70
3:AD:29:THR:HG22	3:AD:30:LYS:H	1.56	0.70
6:AG:150:PHE:HD1	10:AK:55:ARG:HH21	1.39	0.70
23:BB:1854:A:N6	23:BB:1888:G:H1'	2.06	0.70
23:BB:2074:U:H1'	23:BB:2598:A:N3	2.06	0.70
23:BB:600:G:H1'	29:BE:100:MET:HG2	1.74	0.70
25:BC:233:GLY:H	25:BC:241:LYS:NZ	1.90	0.70
26:BD:40:LEU:HA	26:BD:45:TYR:N	2.07	0.70
48:BG:155:PRO:HA	48:BG:170:THR:HG22	1.74	0.70
1:CA:1521:C:H2'	1:CA:1522:U:H6	1.56	0.70
8:CI:81:GLY:O	8:CI:84:ARG:HB2	1.92	0.70
13:CP:4:ILE:O	13:CP:71:VAL:HG11	1.91	0.70
34:D3:54:LEU:HG	34:D3:58:ILE:HD11	1.73	0.70
23:DB:126:A:H5'	36:D2:19:ARG:HG3	1.74	0.70
23:DB:2366:A:H2'	23:DB:2367:G:O4'	1.92	0.70
1:AA:1052:U:H5'	1:AA:1053:G:OP2	1.91	0.70
18:AB:131:LYS:HZ3	18:AB:132:GLU:H	1.40	0.70
2:AC:179:ALA:HA	2:AC:205:GLU:HA	1.73	0.70
16:AS:27:LYS:NZ	16:AS:27:LYS:HB3	2.07	0.70
34:B3:54:LEU:HG	34:B3:58:ILE:HD11	1.72	0.70
23:BB:85:G:OP2	46:BU:6:ARG:HD3	1.92	0.70
23:BB:751:A:H5'	45:BS:90:LYS:HA	1.74	0.70
1:CA:312:C:H2'	1:CA:313:A:H8	1.56	0.70
1:CA:715:A:H2'	1:CA:716:A:C8	2.27	0.70
7:CH:11:THR:HG22	7:CH:14:ARG:HH12	1.57	0.70
8:CI:10:ARG:H	8:CI:80:HIS:CD2	2.10	0.70
11:CL:49:ARG:HG2	11:CL:89:LEU:HD21	1.74	0.70
23:DB:104:A:H2'	23:DB:105:C:H6	1.57	0.70
23:DB:165:A:H2'	23:DB:166:U:H6	1.56	0.70
23:DB:2070:A:H2'	23:DB:2071:A:C8	2.26	0.70
29:DE:34:ALA:HB2	29:DE:96:VAL:HG21	1.74	0.70
28:DP:23:ASP:HA	28:DP:88:ARG:HA	1.72	0.70
1:AA:272:C:H2'	1:AA:273:U:C6	2.26	0.70
4:AE:133:ILE:H	4:AE:133:ILE:HD12	1.57	0.70
8:AI:71:ILE:HD12	8:AI:71:ILE:H	1.55	0.70
16:AS:65:MET:O	16:AS:68:HIS:HB2	1.92	0.70
25:BC:104:LEU:H	25:BC:104:LEU:HD12	1.57	0.70
25:BC:196:ASN:ND2	25:BC:199:HIS:HB2	2.06	0.70
42:BN:97:ILE:HD12	42:BN:98:LEU:H	1.55	0.70
5:CF:67:PRO:O	5:CF:70:VAL:HG22	1.92	0.70
7:CH:11:THR:HA	7:CH:14:ARG:NH1	2.07	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:15:HIS:HA	9:CJ:18:ILE:HG22	1.73	0.70
38:DM:36:VAL:HB	38:DM:127:LYS:O	1.92	0.70
28:DP:63:ILE:HA	28:DP:68:GLY:HA2	1.72	0.70
2:AC:179:ALA:HB1	2:AC:202:PHE:HE1	1.57	0.69
2:AC:82:ASP:O	2:AC:86:LEU:HG	1.92	0.69
7:AH:11:THR:HA	7:AH:14:ARG:NH1	2.07	0.69
9:AJ:77:VAL:HG12	9:AJ:78:GLU:H	1.56	0.69
25:BC:144:GLU:HG3	25:BC:151:GLY:N	2.04	0.69
26:BD:91:THR:HG23	26:BD:92:VAL:H	1.55	0.69
29:BE:60:TRP:O	29:BE:61:ARG:HB2	1.91	0.69
43:BO:27:VAL:HG21	43:BO:40:ILE:HD12	1.74	0.69
49:BR:68:ARG:NH1	49:BR:90:ARG:HD3	2.06	0.69
1:CA:1003:G:H21	1:CA:1005:A:H5'	1.55	0.69
1:CA:1323:G:H2'	1:CA:1324:A:C8	2.27	0.69
1:CA:812:G:O2'	1:CA:813:U:H6	1.74	0.69
23:DB:1515:A:H2'	23:DB:1516:G:O4'	1.92	0.69
23:DB:857:G:C2'	23:DB:858:G:H5'	2.22	0.69
48:DG:42:VAL:HA	48:DG:51:PHE:HA	1.73	0.69
48:DG:84:LYS:HB3	48:DG:132:LEU:O	1.91	0.69
41:DJ:69:ARG:HA	41:DJ:89:PHE:HD1	1.57	0.69
50:DT:92:ASN:HB3	50:DT:93:LEU:HD22	1.74	0.69
1:AA:1162:C:H2'	1:AA:1163:A:C8	2.27	0.69
6:AG:70:PRO:HG3	6:AG:102:TRP:CH2	2.27	0.69
10:AK:28:ASN:ND2	10:AK:46:ALA:HB3	2.06	0.69
12:AM:13:HIS:HB3	12:AM:41:ASP:HA	1.73	0.69
34:B3:49:VAL:HG11	34:B3:54:LEU:HD13	1.74	0.69
1:CA:1278:G:H4'	1:CA:1279:G:C5'	2.22	0.69
1:CA:87:C:C2'	1:CA:88:U:H5'	2.22	0.69
22:DA:43:C:H1'	47:DF:91:ARG:HD2	1.74	0.69
22:DA:49:C:H2'	22:DA:50:A:H8	1.57	0.69
24:DI:105:LEU:HD11	24:DI:139:VAL:HG21	1.72	0.69
39:DX:20:ASN:H	39:DX:20:ASN:HD22	1.40	0.69
1:AA:437:U:H2'	1:AA:438:U:O4'	1.92	0.69
1:AA:406:G:H21	3:AD:115:GLN:HE22	1.40	0.69
12:AM:29:SER:O	12:AM:32:ILE:HG22	1.92	0.69
22:BA:11:C:H3'	22:BA:12:C:H5''	1.73	0.69
23:BB:1459:G:H2'	23:BB:1461:C:C5	2.26	0.69
23:BB:2591:C:H2'	23:BB:2592:G:C8	2.28	0.69
23:BB:2645:G:H5''	23:BB:2732:G:H8	1.57	0.69
23:BB:627:A:H4'	23:BB:628:G:H5'	1.73	0.69
41:BJ:69:ARG:HA	41:BJ:89:PHE:HD1	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:49:ILE:HD13	49:BR:51:VAL:O	1.91	0.69
3:CD:84:ASN:ND2	4:CE:101:GLY:HA3	2.07	0.69
7:CH:103:VAL:HG22	7:CH:124:ILE:HA	1.73	0.69
8:CI:6:TYR:HA	8:CI:18:VAL:O	1.91	0.69
12:CM:44:ILE:H	12:CM:44:ILE:HD12	1.56	0.69
23:DB:1287:A:H3'	23:DB:1288:G:N2	2.07	0.69
25:DC:104:LEU:HD12	25:DC:104:LEU:H	1.57	0.69
2:AC:72:PRO:O	2:AC:76:ILE:HG12	1.92	0.69
23:BB:131:A:H2'	23:BB:132:G:H8	1.56	0.69
23:BB:855:G:H21	52:BW:23:LYS:CG	1.99	0.69
29:BE:134:LEU:HA	29:BE:137:LYS:HB3	1.75	0.69
1:CA:390:U:H2'	1:CA:391:G:C8	2.27	0.69
23:DB:1437:C:H2'	23:DB:1438:U:C6	2.27	0.69
23:DB:558:U:OP1	41:DJ:113:PRO:HG2	1.93	0.69
1:AA:41:G:H2'	1:AA:42:G:C8	2.27	0.69
4:AE:46:GLY:HA3	4:AE:70:MET:HA	1.75	0.69
15:AR:42:ARG:HD2	15:AR:43:ILE:HG23	1.75	0.69
23:BB:1175:A:H3'	23:BB:1176:U:C4'	2.22	0.69
23:BB:2039:U:H2'	23:BB:2040:G:C8	2.27	0.69
29:BE:34:ALA:HB2	29:BE:96:VAL:HG21	1.73	0.69
47:BF:134:GLN:H	47:BF:150:GLY:N	1.91	0.69
40:BH:81:ALA:HB1	40:BH:147:VAL:H	1.56	0.69
35:BV:40:ILE:HD13	35:BV:40:ILE:H	1.57	0.69
1:CA:978:A:H5'	1:CA:1362:A:N6	2.07	0.69
11:CL:86:VAL:HG23	11:CL:89:LEU:HB2	1.74	0.69
14:CQ:74:LEU:HD22	14:CQ:75:VAL:N	2.08	0.69
33:D1:24:LYS:HE2	33:D1:33:LEU:HB2	1.74	0.69
23:DB:1230:A:H2'	23:DB:1231:U:C6	2.27	0.69
23:DB:558:U:H5''	41:DJ:111:LYS:HD3	1.73	0.69
23:DB:858:G:H21	23:DB:2268:A:H3'	1.56	0.69
48:DG:152:ARG:HD3	48:DG:153:PRO:HD2	1.74	0.69
27:DK:54:LYS:H	27:DK:54:LYS:CD	2.05	0.69
43:DO:24:THR:HA	43:DO:42:PRO:HG3	1.75	0.69
35:DV:44:HIS:CE1	35:DV:86:LEU:H	2.11	0.69
51:DZ:77:LYS:HD3	51:DZ:77:LYS:H	1.58	0.69
1:AA:370:C:O2'	1:AA:371:A:H5'	1.92	0.69
2:AC:10:ARG:CZ	2:AC:181:ILE:HD13	2.22	0.69
16:AS:11:ASP:O	16:AS:15:LEU:HB2	1.91	0.69
36:B2:9:VAL:HG13	36:B2:10:LEU:N	2.07	0.69
23:BB:1119:U:OP1	35:BV:83:LYS:HE3	1.92	0.69
26:BD:10:GLY:HA3	26:BD:26:VAL:N	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BO:100:HIS:O	43:BO:104:GLN:HB2	1.92	0.69
43:BO:83:LEU:HD11	43:BO:114:GLY:HA3	1.75	0.69
1:CA:182:A:O2'	1:CA:183:C:H3'	1.92	0.69
8:CI:11:ARG:CZ	8:CI:12:LYS:HB2	2.22	0.69
9:CJ:53:ILE:HG13	21:CN:84:ARG:CZ	2.22	0.69
13:CP:57:ILE:O	13:CP:61:VAL:HG23	1.92	0.69
23:DB:2645:G:H5''	23:DB:2732:G:H8	1.56	0.69
23:DB:856:G:H1'	52:DW:23:LYS:HB3	1.74	0.69
48:DG:84:LYS:HG3	48:DG:131:VAL:HB	1.73	0.69
2:AC:182:ASP:HB3	2:AC:201:ILE:HB	1.74	0.69
12:AM:73:SER:HA	12:AM:76:ILE:HD12	1.74	0.69
23:BB:1871:A:H2'	23:BB:1872:A:C8	2.27	0.69
29:BE:145:ASP:H	29:BE:166:LYS:HB3	1.58	0.69
2:CC:32:LEU:HD11	21:CN:92:ILE:HG12	1.74	0.69
3:CD:12:ARG:NH1	3:CD:12:ARG:HB3	2.08	0.69
3:CD:49:ASP:O	3:CD:53:GLN:HG3	1.92	0.69
8:CI:61:ASP:C	8:CI:62:LEU:HD13	2.13	0.69
23:DB:2147:A:H4'	23:DB:2148:G:C8	2.27	0.69
23:DB:448:U:C5	23:DB:583:G:H1'	2.27	0.69
49:DR:49:ILE:HG12	49:DR:53:PHE:C	2.13	0.69
35:DV:51:GLN:HB2	35:DV:57:TYR:OH	1.91	0.69
1:AA:71:A:N6	1:AA:99:C:H1'	2.02	0.69
23:BB:1937:A:N7	23:BB:1939:U:H2'	2.07	0.69
23:BB:942:G:H2'	23:BB:943:A:O4'	1.93	0.69
29:BE:176:ASP:HB3	29:BE:179:SER:HB2	1.75	0.69
37:BL:103:ILE:H	37:BL:103:ILE:HD12	1.57	0.69
26:BD:186:LEU:HD11	28:BP:3:ILE:HG13	1.75	0.69
45:BS:35:ILE:HG22	45:BS:39:THR:HG21	1.74	0.69
1:CA:1250:A:H4'	8:CI:69:GLY:N	2.07	0.69
1:CA:1296:C:H4'	1:CA:1302:C:H41	1.56	0.69
7:CH:32:LYS:HA	7:CH:35:ILE:HD12	1.75	0.69
8:CI:27:ILE:HD13	8:CI:34:LEU:HB3	1.75	0.69
8:CI:98:ARG:HA	8:CI:103:VAL:HG22	1.74	0.69
17:CT:85:LEU:HD23	17:CT:86:ALA:H	1.57	0.69
23:DB:2502:G:H5'	23:DB:2503:A:H5''	1.74	0.69
40:DH:119:ASN:N	40:DH:119:ASN:HD22	1.90	0.69
23:DB:1080:A:H4'	24:DI:126:ARG:CD	2.22	0.69
37:DL:90:VAL:HB	37:DL:122:VAL:HG12	1.75	0.69
38:DM:34:LYS:O	38:DM:36:VAL:HG23	1.93	0.69
50:DT:60:THR:HB	50:DT:81:LYS:HD2	1.74	0.69
1:AA:1060:U:C4'	9:AJ:54:SER:HB2	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:118:A:H5'	23:BB:119:A:H8	1.56	0.69
23:BB:142:A:H2'	23:BB:143:C:C6	2.28	0.69
23:BB:90:U:H3'	23:BB:91:A:H5''	1.74	0.69
47:BF:163:GLU:HA	47:BF:166:ARG:HD2	1.75	0.69
41:BJ:6:ALA:HB3	41:BJ:45:THR:HG21	1.73	0.69
38:BM:34:LYS:O	38:BM:36:VAL:HG23	1.91	0.69
2:CC:146:LYS:HB2	2:CC:202:PHE:CE2	2.28	0.69
16:CS:30:LEU:HG	16:CS:47:THR:O	1.93	0.69
23:DB:2512:C:H2'	23:DB:2513:A:O4'	1.93	0.69
23:DB:90:U:H3'	23:DB:91:A:H5''	1.73	0.69
41:DJ:45:THR:H	41:DJ:46:PRO:HD3	1.58	0.69
41:DJ:55:ILE:HG22	41:DJ:123:LYS:HB2	1.73	0.69
42:DN:38:LEU:HB3	42:DN:39:PRO:HD3	1.74	0.69
1:AA:1521:C:H2'	1:AA:1522:U:C6	2.28	0.69
18:AB:76:SER:HA	18:AB:92:ASN:HD22	1.58	0.69
5:AF:67:PRO:O	5:AF:70:VAL:HG22	1.93	0.69
23:BB:1035:U:H2'	23:BB:1036:G:H8	1.58	0.69
23:BB:1947:C:H2'	23:BB:1948:G:H8	1.56	0.69
23:BB:2216:G:H2'	23:BB:2217:G:H8	1.58	0.69
26:BD:20:VAL:HG13	27:BK:72:PRO:HB3	1.75	0.69
1:CA:1409:C:H1'	23:DB:1913:A:H62	1.58	0.69
1:CA:41:G:H2'	1:CA:42:G:C8	2.28	0.69
1:CA:844:G:H2'	1:CA:845:A:H8	1.58	0.69
9:CJ:80:THR:HB	9:CJ:82:LYS:HD2	1.75	0.69
22:DA:32:U:H4'	22:DA:52:A:N6	2.08	0.69
49:DR:49:ILE:HD13	49:DR:51:VAL:O	1.92	0.69
35:DV:63:ILE:H	35:DV:63:ILE:HD12	1.58	0.69
1:AA:1060:U:H5''	9:AJ:53:ILE:HG12	1.73	0.69
9:AJ:40:ILE:HB	9:AJ:73:LEU:HB3	1.76	0.69
23:BB:2187:U:H2'	23:BB:2188:U:C6	2.28	0.69
25:BC:183:VAL:HG13	25:BC:185:ALA:H	1.57	0.69
1:CA:1326:U:H2'	1:CA:1327:C:C6	2.27	0.69
6:CG:36:SER:HA	6:CG:39:GLU:OE2	1.92	0.69
10:CK:80:ASN:HA	10:CK:105:ARG:HB3	1.74	0.69
19:CU:39:LYS:N	19:CU:40:PRO:HD2	2.08	0.69
23:DB:546:U:H4'	23:DB:548:G:OP2	1.93	0.69
25:DC:16:VAL:H	25:DC:203:VAL:CG1	2.06	0.69
15:AR:26:ALA:HA	15:AR:29:LYS:HE2	1.75	0.68
16:AS:49:ALA:HA	16:AS:57:VAL:O	1.92	0.68
19:AU:24:LYS:HZ3	19:AU:25:ALA:H	1.37	0.68
23:BB:1060:U:O2	23:BB:1088:A:N7	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2658:C:H5'	48:BG:159:LYS:HZ2	1.57	0.68
23:BB:753:A:H2'	23:BB:754:U:C6	2.27	0.68
25:BC:132:ARG:HH11	25:BC:186:ASP:HB2	1.58	0.68
1:CA:370:C:O2'	1:CA:371:A:H5'	1.93	0.68
3:CD:146:GLU:HA	3:CD:149:LYS:HG2	1.75	0.68
1:CA:8:A:H61	3:CD:53:GLN:HE22	1.41	0.68
10:CK:23:HIS:HB3	10:CK:30:ILE:HG13	1.74	0.68
20:CO:78:TYR:O	20:CO:81:LEU:HB3	1.93	0.68
23:DB:1812:U:H2'	23:DB:1813:G:H8	1.58	0.68
29:DE:134:LEU:HA	29:DE:137:LYS:HB3	1.75	0.68
47:DF:121:PHE:HA	47:DF:127:TYR:HA	1.76	0.68
1:AA:1371:G:O3'	8:AI:70:GLY:HA3	1.92	0.68
3:AD:49:ASP:O	3:AD:53:GLN:HG3	1.94	0.68
13:AP:4:ILE:O	13:AP:71:VAL:HG11	1.92	0.68
44:BQ:45:ALA:O	44:BQ:49:ARG:HB2	1.93	0.68
52:BW:19:ARG:HH21	52:BW:36:ILE:HD11	1.57	0.68
1:CA:73:C:O2'	1:CA:74:A:H5'	1.93	0.68
23:DB:1437:C:H2'	23:DB:1438:U:H6	1.57	0.68
23:DB:1709:U:H2'	23:DB:1710:G:H8	1.58	0.68
23:DB:2039:U:H2'	23:DB:2040:G:C8	2.27	0.68
23:DB:2543:G:H2'	23:DB:2544:G:C8	2.27	0.68
23:DB:597:G:H21	37:DL:12:SER:HA	1.57	0.68
38:DM:19:GLY:HA2	38:DM:97:GLN:HB2	1.75	0.68
45:DS:66:ILE:H	45:DS:66:ILE:HD13	1.58	0.68
1:AA:1461:G:H2'	1:AA:1462:C:H6	1.57	0.68
2:AC:110:LEU:HD12	2:AC:203:LYS:HE2	1.76	0.68
20:AO:67:LEU:HD13	20:AO:88:ARG:HH12	1.56	0.68
23:BB:264:C:H2'	23:BB:265:A:H5'	1.74	0.68
24:BI:10:LEU:HD13	24:BI:12:VAL:HG13	1.74	0.68
41:BJ:45:THR:H	41:BJ:46:PRO:HD3	1.59	0.68
27:BK:54:LYS:H	27:BK:54:LYS:CD	2.05	0.68
27:BK:70:ARG:HB3	27:BK:76:VAL:HG22	1.75	0.68
7:CH:49:LYS:HB3	7:CH:59:GLU:HB2	1.74	0.68
23:DB:2458:G:H8	23:DB:2459:A:H62	1.42	0.68
23:DB:28:A:H61	23:DB:512:G:H1'	1.57	0.68
40:DH:131:SER:CB	40:DH:141:LYS:HA	2.23	0.68
35:DV:40:ILE:HD13	35:DV:40:ILE:H	1.58	0.68
1:AA:1323:G:H2'	1:AA:1324:A:C8	2.27	0.68
1:AA:272:C:H2'	1:AA:273:U:H6	1.59	0.68
1:AA:677:U:H2'	1:AA:678:U:H6	1.58	0.68
10:AK:80:ASN:HA	10:AK:105:ARG:HB3	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:958:A:H61	16:AS:53:GLY:HA3	1.58	0.68
23:BB:2286:G:C8	23:BB:2286:G:H5'	2.28	0.68
23:BB:2543:G:H2'	23:BB:2544:G:C8	2.28	0.68
23:BB:857:G:C2'	23:BB:858:G:H5'	2.23	0.68
49:BR:49:ILE:HG12	49:BR:53:PHE:C	2.14	0.68
4:CE:143:LEU:O	4:CE:146:MET:HG2	1.93	0.68
11:CL:98:ARG:HA	11:CL:103:CYS:SG	2.32	0.68
23:DB:570:G:H2'	23:DB:2030:A:N7	2.09	0.68
23:DB:784:G:C6	25:DC:227:VAL:HG11	2.28	0.68
23:DB:321:U:H1'	29:DE:162:ARG:NH1	2.08	0.68
29:DE:60:TRP:O	29:DE:61:ARG:HB2	1.93	0.68
41:DJ:64:VAL:HG22	41:DJ:68:LYS:HD2	1.74	0.68
43:DO:89:ASP:N	43:DO:115:LEU:HD13	2.07	0.68
35:DV:61:LEU:HD11	35:DV:74:ALA:HB2	1.75	0.68
1:AA:1521:C:H2'	1:AA:1522:U:H6	1.59	0.68
1:AA:182:A:O2'	1:AA:183:C:H3'	1.94	0.68
1:AA:8:A:H61	3:AD:53:GLN:HE22	1.40	0.68
18:AB:198:VAL:HG12	18:AB:200:PRO:HD3	1.73	0.68
18:AB:218:ALA:HA	18:AB:221:ARG:HG2	1.75	0.68
1:AA:392:C:P	13:AP:8:ARG:HH12	2.16	0.68
23:BB:1441:G:H2'	23:BB:1442:U:C6	2.28	0.68
23:BB:1548:A:H2'	23:BB:1549:A:C8	2.29	0.68
23:BB:1553:A:O2'	23:BB:1554:U:H2'	1.93	0.68
23:BB:558:U:OP1	41:BJ:113:PRO:HG2	1.93	0.68
47:BF:34:THR:HA	47:BF:89:THR:HG22	1.75	0.68
44:BQ:4:LYS:NZ	44:BQ:7:VAL:HG22	2.09	0.68
1:CA:1328:C:H5''	12:CM:27:THR:HG21	1.76	0.68
23:DB:1917:U:H2'	23:DB:1918:A:H5'	1.75	0.68
29:DE:176:ASP:HB3	29:DE:179:SER:HB2	1.75	0.68
42:DN:101:GLY:HA2	42:DN:109:PRO:HA	1.74	0.68
1:AA:1349:A:H2'	1:AA:1350:A:O4'	1.93	0.68
18:AB:118:THR:HA	18:AB:121:GLN:HE21	1.58	0.68
8:AI:20:ILE:HG13	8:AI:62:LEU:HD12	1.75	0.68
8:AI:62:LEU:N	8:AI:62:LEU:HD22	2.08	0.68
12:AM:83:GLY:HA2	12:AM:88:LEU:HD21	1.74	0.68
19:AU:39:LYS:N	19:AU:40:PRO:HD2	2.08	0.68
23:BB:1381:G:H2'	23:BB:1382:G:H5'	1.74	0.68
23:BB:2794:C:H2'	23:BB:2795:C:C6	2.28	0.68
23:BB:704:G:H2'	23:BB:726:G:N2	2.08	0.68
47:BF:65:LEU:HD23	47:BF:87:LYS:HD2	1.75	0.68
40:BH:54:LEU:O	40:BH:58:LEU:HB3	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:BK:87:LEU:HD12	27:BK:93:GLN:N	2.08	0.68
38:BM:19:GLY:HA2	38:BM:97:GLN:HB2	1.75	0.68
51:BZ:64:ILE:H	51:BZ:64:ILE:HD12	1.58	0.68
23:DB:138:U:O3'	23:DB:139:U:H3'	1.92	0.68
23:DB:483:A:H5'	46:DU:44:HIS:O	1.94	0.68
26:DD:20:VAL:HG13	27:DK:72:PRO:HB3	1.75	0.68
27:DK:25:LEU:HG	27:DK:39:ILE:HA	1.75	0.68
50:DT:2:ILE:HD11	50:DT:3:ARG:CZ	2.24	0.68
1:AA:1308:U:H3'	12:AM:97:ARG:NH1	2.08	0.68
5:AF:2:ARG:HG3	5:AF:92:THR:OG1	1.94	0.68
1:AA:1179:A:H4'	8:AI:104:THR:HA	1.76	0.68
9:AJ:55:PRO:HA	21:AN:80:ARG:NH2	2.09	0.68
22:BA:91:C:H2'	22:BA:92:C:H6	1.59	0.68
23:BB:1041:G:H2'	23:BB:1042:G:H8	1.58	0.68
47:BF:66:ILE:HD12	47:BF:85:GLY:O	1.94	0.68
40:BH:131:SER:HB3	40:BH:140:ALA:O	1.94	0.68
52:BW:23:LYS:HD2	52:BW:24:ARG:H	1.58	0.68
1:CA:1448:C:H2'	1:CA:1449:C:H6	1.59	0.68
4:CE:156:ARG:NH1	7:CH:42:GLU:HB3	2.09	0.68
23:DB:2328:A:H2'	23:DB:2329:U:H6	1.59	0.68
23:DB:855:G:N2	52:DW:23:LYS:HG2	2.07	0.68
50:DT:54:GLU:HB3	50:DT:88:LYS:HB2	1.74	0.68
46:DU:40:LEU:H	46:DU:40:LEU:HD12	1.58	0.68
1:AA:16:A:O2'	1:AA:17:U:H5'	1.94	0.68
1:AA:764:C:H3'	1:AA:765:G:N2	2.08	0.68
18:AB:185:ILE:HA	18:AB:199:ILE:O	1.94	0.68
2:AC:154:GLY:HA3	2:AC:162:ALA:HB1	1.74	0.68
7:AH:32:LYS:HA	7:AH:35:ILE:HD12	1.75	0.68
12:AM:79:LEU:HA	12:AM:82:LEU:HD12	1.76	0.68
1:AA:1308:U:OP1	12:AM:95:PRO:HA	1.93	0.68
23:BB:2389:G:H5''	23:BB:2390:U:H5'	1.75	0.68
26:BD:54:ALA:HA	26:BD:76:GLY:HA2	1.76	0.68
26:BD:69:ALA:HA	26:BD:73:VAL:HB	1.76	0.68
48:BG:84:LYS:HB2	48:BG:132:LEU:HD23	1.75	0.68
43:BO:24:THR:HA	43:BO:42:PRO:HG3	1.75	0.68
3:CD:153:ARG:HG3	3:CD:154:VAL:H	1.58	0.68
7:CH:54:THR:HG23	7:CH:55:LYS:H	1.58	0.68
7:CH:63:LYS:HG2	7:CH:70:VAL:HG21	1.76	0.68
8:CI:11:ARG:HB2	8:CI:106:ASP:HB2	1.76	0.68
8:CI:87:MET:HB2	8:CI:94:ARG:NH2	2.09	0.68
23:DB:753:A:H2'	23:DB:754:U:C6	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:946:C:H2'	23:DB:947:A:H8	1.59	0.68
26:DD:106:LYS:HB3	26:DD:206:ALA:N	2.08	0.68
24:DI:73:PRO:HG2	24:DI:78:LEU:HD21	1.75	0.68
41:DJ:3:THR:HG21	44:DQ:60:TRP:HE1	1.59	0.68
23:DB:974:G:OP2	49:DR:78:ARG:HD3	1.94	0.68
52:DW:19:ARG:HH21	52:DW:36:ILE:HD11	1.57	0.68
2:AC:52:SER:HB3	2:AC:114:LEU:HD21	1.75	0.68
23:BB:2070:A:H2'	23:BB:2071:A:C8	2.29	0.68
23:BB:2819:G:H2'	23:BB:2821:A:N7	2.08	0.68
42:BN:47:VAL:C	42:BN:50:PRO:HD2	2.14	0.68
43:BO:89:ASP:N	43:BO:115:LEU:HD13	2.08	0.68
35:BV:61:LEU:HD11	35:BV:74:ALA:HB2	1.76	0.68
2:CC:120:THR:HG23	2:CC:197:VAL:HG21	1.74	0.68
48:DG:148:ARG:HD3	48:DG:152:ARG:NH1	2.09	0.68
48:DG:84:LYS:HB2	48:DG:132:LEU:HD23	1.76	0.68
40:DH:87:GLU:OE2	40:DH:89:LYS:HD3	1.93	0.68
27:DK:10:VAL:HG13	27:DK:86:LEU:HD21	1.75	0.68
38:DM:126:ILE:H	38:DM:126:ILE:HD12	1.58	0.68
28:DP:97:TYR:O	28:DP:100:ARG:HB2	1.94	0.68
1:AA:844:G:H2'	1:AA:845:A:H8	1.59	0.68
2:AC:179:ALA:HB3	2:AC:181:ILE:HD11	1.76	0.68
23:BB:1381:G:C2'	23:BB:1382:G:H5'	2.24	0.68
23:BB:1437:C:H2'	23:BB:1438:U:C6	2.29	0.68
23:BB:1515:A:H2'	23:BB:1516:G:O4'	1.94	0.68
1:CA:71:A:O2'	1:CA:72:A:H5'	1.94	0.68
18:CB:75:ALA:C	18:CB:79:VAL:HG23	2.15	0.68
10:CK:34:THR:HA	10:CK:41:LEU:HG	1.75	0.68
23:DB:414:C:H2'	23:DB:415:A:C8	2.28	0.68
47:DF:109:ARG:HB2	47:DF:135:ILE:HD12	1.76	0.68
1:AA:1081:A:H2'	1:AA:1082:A:H8	1.59	0.67
1:AA:8:A:N6	3:AD:205:LYS:HB2	2.08	0.67
17:AT:38:ILE:HD11	17:AT:82:ILE:HG22	1.77	0.67
23:BB:2737:G:H2'	23:BB:2738:A:C8	2.29	0.67
48:BG:42:VAL:HA	48:BG:51:PHE:HA	1.76	0.67
41:BJ:64:VAL:HG22	41:BJ:68:LYS:HD2	1.75	0.67
27:BK:118:LEU:C	27:BK:120:PRO:HD2	2.13	0.67
45:BS:66:ILE:H	45:BS:66:ILE:HD13	1.57	0.67
1:CA:543:U:H2'	1:CA:544:G:H8	1.58	0.67
1:CA:900:A:H2'	1:CA:901:A:C8	2.29	0.67
1:CA:973:G:H3'	1:CA:974:A:H5''	1.75	0.67
23:DB:1947:C:H2'	23:DB:1948:G:H8	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2187:U:H2'	23:DB:2188:U:C6	2.28	0.67
23:DB:2591:C:H2'	23:DB:2592:G:C8	2.29	0.67
23:DB:2841:C:H2'	23:DB:2842:G:H8	1.59	0.67
1:AA:1118:U:H5'	8:AI:10:ARG:NH1	2.09	0.67
8:AI:78:ILE:HG22	8:AI:82:ILE:HD11	1.76	0.67
36:B2:13:ASN:O	36:B2:17:GLY:N	2.27	0.67
23:BB:1438:U:H2'	23:BB:1439:A:O4'	1.94	0.67
23:BB:165:A:H2'	23:BB:166:U:H6	1.59	0.67
26:BD:16:THR:HB	26:BD:18:ASP:OD1	1.93	0.67
23:BB:2658:C:H5'	48:BG:159:LYS:NZ	2.08	0.67
50:BT:92:ASN:HB3	50:BT:93:LEU:HD22	1.75	0.67
39:BX:20:ASN:H	39:BX:20:ASN:HD22	1.43	0.67
1:CA:392:C:P	13:CP:8:ARG:HH12	2.16	0.67
1:CA:8:A:N6	3:CD:205:LYS:HB2	2.09	0.67
1:CA:658:C:O3'	20:CO:8:THR:HG21	1.94	0.67
16:CS:29:PRO:HA	16:CS:47:THR:HB	1.76	0.67
36:D2:9:VAL:HG13	36:D2:10:LEU:N	2.08	0.67
23:DB:1060:U:O2	23:DB:1088:A:N7	2.27	0.67
23:DB:1283:G:H22	23:DB:1286:A:H5'	1.59	0.67
23:DB:2841:C:H2'	23:DB:2842:G:C8	2.29	0.67
26:DD:40:LEU:HA	26:DD:45:TYR:N	2.09	0.67
24:DI:1:ALA:HB1	24:DI:2:LYS:HD2	1.77	0.67
37:DL:79:LEU:HB2	37:DL:113:ALA:H	1.59	0.67
43:DO:68:LYS:HA	43:DO:102:ARG:HG2	1.76	0.67
1:AA:524:G:H2'	1:AA:525:C:C6	2.29	0.67
1:AA:989:U:H2'	1:AA:990:C:O4'	1.94	0.67
17:AT:85:LEU:HD23	17:AT:86:ALA:H	1.59	0.67
23:BB:1230:A:H2'	23:BB:1231:U:C6	2.29	0.67
23:BB:1252:G:H1'	44:BQ:32:ARG:HH22	1.59	0.67
23:BB:1283:G:H22	23:BB:1286:A:H5'	1.59	0.67
35:BV:38:LEU:HG	35:BV:40:ILE:HD12	1.75	0.67
1:CA:1268:G:H2'	1:CA:1269:A:C8	2.30	0.67
1:CA:250:A:H1'	1:CA:252:U:C6	2.29	0.67
1:CA:328:C:H4'	1:CA:329:A:H5''	1.76	0.67
3:CD:197:HIS:O	3:CD:200:VAL:HG22	1.95	0.67
9:CJ:6:ILE:HB	9:CJ:76:ILE:CG1	2.24	0.67
10:CK:75:GLU:H	10:CK:75:GLU:CD	1.97	0.67
23:DB:1487:U:H2'	23:DB:1488:C:C6	2.29	0.67
23:DB:2748:A:H4'	48:DG:3:VAL:HG21	1.76	0.67
37:DL:103:ILE:H	37:DL:103:ILE:HD12	1.57	0.67
1:AA:86:G:H21	1:AA:87:C:N4	1.91	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AS:35:ARG:HB2	16:AS:71:GLY:HA2	1.75	0.67
23:BB:2867:G:N3	23:BB:2867:G:H2'	2.09	0.67
23:BB:67:U:H2'	23:BB:68:G:H8	1.58	0.67
25:BC:68:ARG:HH21	25:BC:190:THR:HG22	1.59	0.67
26:BD:106:LYS:HB3	26:BD:206:ALA:N	2.09	0.67
35:BV:5:ASN:HA	35:BV:64:VAL:HB	1.77	0.67
1:CA:1391:U:H2'	1:CA:1392:G:C8	2.30	0.67
1:CA:803:G:H2'	1:CA:804:U:C6	2.30	0.67
2:CC:76:ILE:HD11	2:CC:102:ILE:HG21	1.76	0.67
10:CK:17:ASP:HA	10:CK:80:ASN:O	1.94	0.67
21:CN:31:SER:HA	21:CN:44:VAL:HB	1.76	0.67
23:DB:1124:G:H1'	32:D4:38:GLY:OXT	1.94	0.67
23:DB:1729:U:H5''	23:DB:1730:C:O2	1.95	0.67
23:DB:2704:C:H2'	23:DB:2705:A:O4'	1.95	0.67
23:DB:560:C:H3'	23:DB:561:G:C8	2.30	0.67
23:DB:67:U:H2'	23:DB:68:G:H8	1.57	0.67
27:DK:112:PHE:O	27:DK:115:ILE:HG22	1.93	0.67
44:DQ:57:ARG:NH1	44:DQ:61:ILE:HD11	2.09	0.67
49:DR:39:LEU:HB3	49:DR:49:ILE:HD11	1.75	0.67
52:DW:35:ILE:HA	52:DW:57:THR:HG23	1.74	0.67
1:AA:328:C:H4'	1:AA:329:A:H5''	1.76	0.67
1:AA:845:A:H5''	1:AA:846:G:C8	2.30	0.67
21:AN:50:LEU:N	21:AN:51:PRO:HD2	2.05	0.67
23:BB:1484:U:H2'	23:BB:1485:U:C6	2.29	0.67
23:BB:1760:C:H2'	23:BB:1761:C:O4'	1.95	0.67
23:BB:184:C:H2'	23:BB:185:G:C8	2.29	0.67
23:BB:2086:U:H2'	23:BB:2087:G:C8	2.30	0.67
23:BB:477:A:H2'	23:BB:478:A:H8	1.59	0.67
44:BQ:89:ILE:HG21	49:BR:11:GLN:HE22	1.59	0.67
30:BY:3:THR:HB	30:BY:36:GLU:HG3	1.74	0.67
1:CA:1086:U:H3	1:CA:1099:G:H22	1.42	0.67
18:CB:63:LYS:HA	18:CB:224:ARG:HH11	1.60	0.67
2:CC:5:HIS:NE2	2:CC:7:ASN:HB3	2.10	0.67
8:CI:17:ARG:O	8:CI:64:ILE:HG23	1.94	0.67
17:CT:43:LYS:HD3	17:CT:43:LYS:H	1.59	0.67
31:D0:9:ARG:O	31:D0:12:ARG:HB3	1.94	0.67
23:DB:1657:U:O2'	23:DB:1658:C:H5'	1.95	0.67
23:DB:2086:U:H2'	23:DB:2087:G:C8	2.30	0.67
23:DB:2094:A:H2'	23:DB:2095:A:C8	2.30	0.67
23:DB:2341:G:H2'	23:DB:2342:C:C6	2.30	0.67
23:DB:2789:C:H3'	23:DB:2893:A:H62	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DM:33:LEU:HD22	38:DM:128:THR:HB	1.77	0.67
42:DN:48:VAL:HA	42:DN:51:LEU:HD22	1.77	0.67
1:AA:724:G:H2'	1:AA:725:G:H8	1.58	0.67
18:AB:117:GLU:O	18:AB:121:GLN:HB2	1.95	0.67
23:BB:27:G:N2	23:BB:512:G:H2'	2.08	0.67
23:BB:1076:C:H4'	24:BI:94:LYS:HE3	1.77	0.67
28:BP:97:TYR:O	28:BP:100:ARG:HB2	1.95	0.67
1:CA:1238:A:H5'	1:CA:1336:C:H41	1.56	0.67
1:CA:1461:G:H2'	1:CA:1462:C:C6	2.29	0.67
1:CA:16:A:O2'	1:CA:17:U:H5'	1.95	0.67
1:CA:524:G:H2'	1:CA:525:C:C6	2.29	0.67
16:CS:18:VAL:O	16:CS:22:VAL:HG23	1.94	0.67
23:DB:1047:G:H1'	23:DB:1110:G:N2	2.09	0.67
23:DB:165:A:H2'	23:DB:166:U:C6	2.30	0.67
23:DB:191:A:H2'	23:DB:192:C:C6	2.30	0.67
23:DB:974:G:H1'	23:DB:975:A:C8	2.30	0.67
27:DK:70:ARG:HB3	27:DK:76:VAL:HG22	1.76	0.67
44:DQ:30:VAL:HG13	44:DQ:31:TYR:N	2.09	0.67
44:DQ:91:ARG:HD3	49:DR:11:GLN:HB2	1.76	0.67
1:AA:958:A:N6	16:AS:53:GLY:HA3	2.09	0.67
12:AM:106:ARG:HE	12:AM:112:ARG:HG2	1.58	0.67
23:BB:2046:G:H1'	31:B0:18:HIS:CD2	2.30	0.67
23:BB:2291:U:H2'	23:BB:2292:U:C6	2.29	0.67
23:BB:365:U:H2'	23:BB:366:C:C6	2.30	0.67
26:BD:55:LYS:HZ2	26:BD:60:VAL:HG13	1.58	0.67
48:BG:152:ARG:HD3	48:BG:153:PRO:HD2	1.76	0.67
48:BG:157:LYS:HG2	48:BG:159:LYS:HG3	1.77	0.67
48:BG:89:VAL:HG12	48:BG:90:GLY:H	1.60	0.67
38:BM:36:VAL:HB	38:BM:127:LYS:O	1.95	0.67
23:DB:1080:A:H2'	23:DB:1081:U:H6	1.57	0.67
23:DB:1441:G:H2'	23:DB:1442:U:C6	2.30	0.67
23:DB:228:C:H4'	23:DB:229:C:H5'	1.76	0.67
29:DE:147:LEU:HD12	29:DE:149:ILE:HB	1.76	0.67
47:DF:65:LEU:HD23	47:DF:87:LYS:HD2	1.75	0.67
41:DJ:124:VAL:HG23	41:DJ:125:TYR:H	1.59	0.67
38:DM:82:MET:HE3	38:DM:83:GLY:H	1.60	0.67
45:DS:35:ILE:HG22	45:DS:39:THR:HG21	1.76	0.67
2:AC:40:GLN:HG3	2:AC:41:TYR:N	2.09	0.67
4:AE:143:LEU:O	4:AE:146:MET:HG2	1.95	0.67
8:AI:41:GLU:N	8:AI:44:ARG:HH11	1.93	0.67
12:AM:53:ASP:HA	12:AM:56:ARG:NH2	2.10	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B1:24:LYS:HE2	33:B1:33:LEU:HB2	1.77	0.67
29:BE:147:LEU:HD12	29:BE:149:ILE:HB	1.76	0.67
44:BQ:30:VAL:HG13	44:BQ:31:TYR:N	2.10	0.67
1:CA:17:U:H2'	1:CA:18:C:H6	1.57	0.67
1:CA:817:C:H1'	1:CA:819:A:H5'	1.77	0.67
4:CE:46:GLY:HA3	4:CE:70:MET:HA	1.77	0.67
9:CJ:12:ALA:HB2	9:CJ:96:VAL:HG12	1.76	0.67
16:CS:39:ILE:HG13	16:CS:68:HIS:O	1.95	0.67
23:DB:1139:G:O2'	23:DB:1140:C:H5'	1.95	0.67
23:DB:1176:U:H3'	23:DB:1177:G:C8	2.29	0.67
23:DB:1381:G:C2'	23:DB:1382:G:H5'	2.24	0.67
23:DB:1484:U:H2'	23:DB:1485:U:C6	2.28	0.67
23:DB:2143:C:H3'	23:DB:2144:G:O4'	1.94	0.67
23:DB:454:A:H3'	23:DB:455:C:H5'	1.76	0.67
23:DB:2307:G:O6	47:DF:84:ILE:HD11	1.95	0.67
52:DW:37:VAL:HG13	52:DW:55:ASP:O	1.95	0.67
18:AB:46:VAL:HA	18:AB:49:PHE:CD2	2.30	0.67
2:AC:133:MET:CE	2:AC:152:VAL:HG21	2.25	0.67
1:AA:532:A:N6	2:AC:191:THR:HB	2.10	0.67
6:AG:149:ALA:N	10:AK:55:ARG:HH22	1.92	0.67
11:AL:13:ARG:O	11:AL:14:LYS:HB3	1.94	0.67
23:BB:215:G:H4'	23:BB:216:A:H4'	1.77	0.67
23:BB:832:U:H2'	23:BB:833:A:C8	2.29	0.67
47:BF:102:LEU:HA	47:BF:106:ALA:HB2	1.75	0.67
38:BM:126:ILE:HD12	38:BM:126:ILE:H	1.60	0.67
1:CA:190:A:H2'	1:CA:191:G:O4'	1.94	0.67
6:CG:57:GLU:HA	6:CG:60:ALA:HB3	1.76	0.67
23:DB:910:A:H2'	23:DB:911:A:C8	2.29	0.67
25:DC:93:VAL:HG13	25:DC:94:LEU:N	2.08	0.67
46:DU:85:ARG:HD3	46:DU:86:PHE:H	1.60	0.67
22:BA:60:C:H2'	22:BA:61:G:H8	1.59	0.67
23:BB:1657:U:O2'	23:BB:1658:C:H5'	1.95	0.67
23:BB:2461:A:H2'	23:BB:2462:C:C6	2.30	0.67
23:BB:2674:G:H4'	27:BK:30:ARG:HG3	1.77	0.67
23:BB:547:A:H3'	23:BB:548:G:O4'	1.96	0.67
23:BB:83:A:H61	23:BB:101:A:H5'	1.59	0.67
25:BC:143:VAL:HB	25:BC:153:LEU:HB2	1.75	0.67
25:BC:75:ALA:HB2	25:BC:95:TYR:HA	1.76	0.67
47:BF:141:ASP:O	47:BF:145:VAL:HG13	1.95	0.67
40:BH:116:ARG:NH1	40:BH:133:GLN:HB2	2.10	0.67
41:BJ:11:VAL:HG11	41:BJ:13:ARG:HE	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BQ:77:LYS:HB3	44:BQ:116:LEU:HD21	1.77	0.67
45:BS:66:ILE:HA	45:BS:69:LEU:HD13	1.75	0.67
1:CA:1071:C:H2'	1:CA:1072:G:H8	1.60	0.67
1:CA:301:G:H2'	1:CA:302:G:C8	2.29	0.67
20:CO:36:ILE:HD12	20:CO:60:VAL:HG23	1.75	0.67
13:CP:6:LEU:HD11	13:CP:71:VAL:HB	1.77	0.67
23:DB:2021:C:OP1	31:D0:8:THR:HG21	1.95	0.67
40:DH:27:ARG:NH1	51:DZ:60:ASP:HA	2.10	0.67
27:DK:66:LYS:HA	27:DK:79:PHE:O	1.94	0.67
28:DP:89:GLY:HA2	28:DP:112:ARG:N	2.09	0.67
45:DS:26:GLY:N	45:DS:71:VAL:HG13	2.09	0.67
1:AA:202:G:H21	1:AA:465:A:H61	1.42	0.66
1:AA:859:G:H2'	1:AA:860:A:H8	1.61	0.66
23:BB:543:G:H2'	23:BB:544:C:C5'	2.25	0.66
23:BB:932:U:H5'	23:BB:933:A:OP1	1.95	0.66
44:BQ:89:ILE:HG22	44:BQ:91:ARG:H	1.60	0.66
50:BT:69:ARG:HG2	50:BT:73:ARG:C	2.15	0.66
1:CA:1244:G:H2'	1:CA:1245:C:C6	2.30	0.66
3:CD:2:ARG:NH1	3:CD:114:ARG:HD3	2.09	0.66
12:CM:29:SER:O	12:CM:33:LEU:HG	1.95	0.66
9:CJ:52:LEU:HB2	21:CN:80:ARG:HD3	1.77	0.66
14:CQ:58:VAL:HB	14:CQ:74:LEU:HD23	1.77	0.66
23:DB:1438:U:H2'	23:DB:1439:A:O4'	1.95	0.66
23:DB:1480:C:H2'	23:DB:1481:U:O4'	1.95	0.66
23:DB:2819:G:H2'	23:DB:2821:A:N7	2.09	0.66
23:DB:2886:A:H62	31:D0:39:ARG:CZ	2.07	0.66
26:DD:40:LEU:HD23	26:DD:46:ARG:HG2	1.77	0.66
29:DE:33:VAL:O	29:DE:36:ALA:HB3	1.96	0.66
48:DG:91:VAL:HG23	48:DG:92:GLY:H	1.61	0.66
41:DJ:109:LEU:HD13	41:DJ:119:PHE:HB2	1.77	0.66
41:DJ:64:VAL:O	41:DJ:65:THR:HG22	1.95	0.66
37:DL:143:GLU:CG	37:DL:144:GLU:H	2.07	0.66
1:AA:1163:A:H2'	1:AA:1164:G:H8	1.61	0.66
3:AD:153:ARG:HG3	3:AD:154:VAL:H	1.59	0.66
5:AF:6:ILE:HG23	5:AF:62:MET:HB3	1.76	0.66
6:AG:67:ASN:ND2	6:AG:127:ALA:HA	2.11	0.66
13:AP:6:LEU:HD11	13:AP:71:VAL:HB	1.77	0.66
23:BB:2704:C:H2'	23:BB:2705:A:O4'	1.95	0.66
25:BC:93:VAL:HG13	25:BC:94:LEU:N	2.10	0.66
40:BH:4:ILE:HB	40:BH:37:VAL:CG1	2.25	0.66
1:CA:272:C:H2'	1:CA:273:U:C6	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:926:G:N2	1:CA:1505:G:H2'	2.10	0.66
3:CD:84:ASN:ND2	3:CD:86:GLY:H	1.93	0.66
9:CJ:52:LEU:HA	9:CJ:62:ARG:HA	1.76	0.66
11:CL:49:ARG:HB2	11:CL:89:LEU:HD11	1.76	0.66
15:CR:42:ARG:HD2	15:CR:43:ILE:HG23	1.77	0.66
22:DA:25:U:H4'	22:DA:26:C:H5''	1.77	0.66
23:DB:1760:C:H2'	23:DB:1761:C:O4'	1.95	0.66
23:DB:2548:U:H1'	27:DK:23:LYS:HZ1	1.57	0.66
23:DB:2867:G:N3	23:DB:2867:G:H2'	2.10	0.66
23:DB:38:A:N3	29:DE:43:THR:HB	2.10	0.66
44:DQ:77:LYS:HB3	44:DQ:116:LEU:HD21	1.76	0.66
45:DS:66:ILE:HA	45:DS:69:LEU:HD13	1.77	0.66
35:DV:5:ASN:HA	35:DV:64:VAL:HB	1.77	0.66
1:AA:1125:U:C2'	1:AA:1126:U:H5''	2.25	0.66
1:AA:1194:U:H2'	1:AA:1195:C:C6	2.29	0.66
1:AA:1298:U:H4'	1:AA:1299:A:O5'	1.96	0.66
1:AA:859:G:H2'	1:AA:860:A:C8	2.31	0.66
18:AB:30:ILE:HG23	18:AB:38:HIS:HB3	1.77	0.66
7:AH:63:LYS:HG2	7:AH:70:VAL:HG21	1.77	0.66
23:BB:1487:U:H2'	23:BB:1488:C:C6	2.30	0.66
23:BB:454:A:H3'	23:BB:455:C:H5''	1.77	0.66
50:BT:34:VAL:HG11	50:BT:43:ILE:HD11	1.77	0.66
1:CA:193:C:H2'	1:CA:194:C:C6	2.31	0.66
1:CA:764:C:H3'	1:CA:765:G:N2	2.10	0.66
18:CB:186:VAL:O	18:CB:200:PRO:HA	1.95	0.66
5:CF:72:ASP:HA	5:CF:75:GLU:OE1	1.95	0.66
23:DB:1935:G:H1'	23:DB:1964:G:N2	2.10	0.66
23:DB:2286:G:H5'	23:DB:2286:G:C8	2.30	0.66
23:DB:2893:A:H5''	23:DB:2894:G:H5'	1.78	0.66
23:DB:2:G:H2'	23:DB:3:U:H6	1.59	0.66
25:DC:75:ALA:HB2	25:DC:95:TYR:HA	1.76	0.66
27:DK:73:ASP:O	28:DP:74:GLN:HG3	1.95	0.66
7:AH:54:THR:HG23	7:AH:55:LYS:H	1.60	0.66
12:AM:106:ARG:HH11	12:AM:109:LYS:HD2	1.60	0.66
23:BB:1594:U:H2'	23:BB:1595:C:C6	2.31	0.66
23:BB:1704:C:H2'	23:BB:1705:A:C8	2.30	0.66
23:BB:2586:U:H2'	23:BB:2587:A:C8	2.29	0.66
26:BD:148:GLN:CB	26:BD:152:PRO:HG2	2.26	0.66
23:BB:588:U:H1'	29:BE:85:PHE:CD2	2.31	0.66
47:BF:111:ARG:HH22	47:BF:113:PHE:HB2	1.59	0.66
27:BK:66:LYS:HA	27:BK:79:PHE:O	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BU:13:LEU:HA	46:BU:18:LYS:HE3	1.78	0.66
1:CA:724:G:H2'	1:CA:725:G:H8	1.59	0.66
18:CB:98:GLY:C	18:CB:106:VAL:HG21	2.16	0.66
1:CA:1108:G:H5'	2:CC:175:HIS:ND1	2.10	0.66
20:CO:35:GLN:O	20:CO:39:LEU:HD13	1.94	0.66
32:D4:16:ILE:HA	32:D4:25:VAL:HG22	1.78	0.66
32:D4:22:VAL:HG11	32:D4:36:ARG:HD2	1.77	0.66
23:DB:2389:G:H5''	23:DB:2390:U:H5'	1.77	0.66
26:DD:69:ALA:HA	26:DD:73:VAL:HB	1.76	0.66
29:DE:145:ASP:H	29:DE:166:LYS:HB3	1.59	0.66
48:DG:30:GLY:CA	48:DG:78:VAL:HA	2.24	0.66
44:DQ:45:ALA:O	44:DQ:49:ARG:HB2	1.95	0.66
51:DZ:54:LYS:HA	51:DZ:57:ARG:HD3	1.78	0.66
1:AA:880:C:H2'	1:AA:881:G:H8	1.60	0.66
1:AA:1118:U:H5'	8:AI:10:ARG:HH11	1.60	0.66
10:AK:16:SER:HA	10:AK:78:ILE:HA	1.78	0.66
20:AO:26:GLU:HG2	20:AO:81:LEU:HD12	1.77	0.66
13:AP:1:MET:HE1	13:AP:66:THR:OG1	1.96	0.66
14:AQ:58:VAL:HB	14:AQ:74:LEU:HD23	1.76	0.66
32:B4:5:ALA:HA	32:B4:37:GLN:NE2	2.10	0.66
23:BB:1935:G:H1'	23:BB:1964:G:N2	2.11	0.66
23:BB:2097:A:H2'	23:BB:2098:U:C6	2.30	0.66
23:BB:2267:A:C8	23:BB:2267:A:C3'	2.79	0.66
23:BB:878:A:N3	23:BB:878:A:H2'	2.10	0.66
25:BC:16:VAL:H	25:BC:203:VAL:CG1	2.09	0.66
41:BJ:64:VAL:O	41:BJ:65:THR:HG22	1.96	0.66
43:BO:68:LYS:HA	43:BO:102:ARG:HG2	1.77	0.66
44:BQ:105:PHE:HA	44:BQ:108:LEU:HB2	1.78	0.66
35:BV:44:HIS:CE1	35:BV:86:LEU:H	2.14	0.66
35:BV:63:ILE:HD12	35:BV:63:ILE:H	1.61	0.66
21:CN:65:GLN:H	21:CN:65:GLN:HE21	1.40	0.66
13:CP:1:MET:HE1	13:CP:66:THR:OG1	1.95	0.66
23:DB:2425:A:H5'	23:DB:2427:C:O4'	1.95	0.66
48:DG:153:PRO:HB3	48:DG:158:GLY:HA2	1.77	0.66
46:DU:13:LEU:HA	46:DU:18:LYS:HE3	1.77	0.66
6:AG:72:VAL:O	6:AG:140:VAL:HG12	1.96	0.66
10:AK:34:THR:HA	10:AK:41:LEU:HG	1.77	0.66
36:B2:9:VAL:HG13	36:B2:10:LEU:H	1.59	0.66
23:BB:328:U:H4'	46:BU:65:GLN:NE2	2.10	0.66
23:BB:28:A:N6	23:BB:512:G:H1'	2.11	0.66
26:BD:116:LYS:HB3	26:BD:118:PHE:CZ	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:58:LYS:HZ2	29:BE:58:LYS:H	1.42	0.66
23:BB:196:A:H5''	37:BL:47:ARG:HH12	1.59	0.66
50:BT:5:GLU:O	50:BT:9:LYS:HE3	1.96	0.66
1:CA:1230:C:H2'	1:CA:1231:G:H8	1.61	0.66
1:CA:518:C:H2'	1:CA:530:G:C8	2.30	0.66
1:CA:958:A:N6	16:CS:76:THR:HG22	2.10	0.66
6:CG:14:ASP:HB3	6:CG:18:GLY:N	2.11	0.66
8:CI:56:MET:C	8:CI:58:GLU:H	1.99	0.66
12:CM:95:PRO:N	12:CM:108:ARG:HG2	2.10	0.66
23:DB:1024:G:C3'	23:DB:1025:G:H5''	2.26	0.66
23:DB:2150:C:H2'	23:DB:2151:U:C6	2.30	0.66
23:DB:2469:A:H4'	38:DM:55:ARG:NE	2.10	0.66
23:DB:27:G:N2	23:DB:512:G:H2'	2.08	0.66
23:DB:634:C:H2'	23:DB:635:C:H6	1.61	0.66
1:AA:1105:A:H2'	1:AA:1106:G:H8	1.60	0.66
1:AA:1169:A:H2'	1:AA:1170:A:C8	2.30	0.66
1:AA:215:C:H2'	1:AA:216:U:C6	2.30	0.66
1:AA:429:U:H1'	1:AA:430:A:H5''	1.78	0.66
1:AA:715:A:H2'	1:AA:716:A:H8	1.61	0.66
18:AB:13:VAL:O	18:AB:14:HIS:HB3	1.95	0.66
4:AE:113:VAL:HG23	4:AE:114:LEU:N	2.11	0.66
26:BD:33:ARG:HE	26:BD:51:THR:HB	1.60	0.66
23:BB:674:G:O3'	29:BE:60:TRP:CZ2	2.49	0.66
48:BG:17:LYS:HE2	48:BG:19:ASN:ND2	2.11	0.66
37:BL:143:GLU:CG	37:BL:144:GLU:H	2.09	0.66
38:BM:78:LEU:HD12	38:BM:78:LEU:H	1.61	0.66
35:BV:72:VAL:HB	35:BV:92:VAL:O	1.95	0.66
52:BW:23:LYS:HD2	52:BW:24:ARG:HB3	1.77	0.66
39:BX:13:GLU:HB2	39:BX:57:LEU:HD23	1.77	0.66
1:CA:423:G:H2'	1:CA:424:G:O4'	1.96	0.66
18:CB:137:THR:HA	18:CB:140:LEU:HD13	1.78	0.66
33:D1:9:LYS:H	33:D1:9:LYS:HD3	1.60	0.66
22:DA:60:C:H2'	22:DA:61:G:H8	1.61	0.66
23:DB:1821:A:H2'	23:DB:1822:C:C6	2.31	0.66
23:DB:28:A:N6	23:DB:512:G:H1'	2.11	0.66
23:DB:455:C:N3	23:DB:472:A:H2'	2.11	0.66
47:DF:111:ARG:HH22	47:DF:113:PHE:HB2	1.61	0.66
46:DU:78:LYS:HD3	46:DU:79:ALA:H	1.61	0.66
23:DB:96:C:H4'	39:DX:41:HIS:ND1	2.11	0.66
30:DY:7:THR:HG23	30:DY:34:THR:OG1	1.95	0.66
1:AA:1486:G:H2'	1:AA:1487:G:O4'	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:312:C:H2'	1:AA:313:A:H8	1.59	0.66
2:AC:129:PHE:CG	2:AC:130:ARG:N	2.63	0.66
1:AA:1057:G:H5''	2:AC:153:SER:CB	2.25	0.66
2:AC:181:ILE:N	2:AC:181:ILE:HD12	2.09	0.66
10:AK:124:LYS:HA	19:AU:34:ARG:HB3	1.78	0.66
12:AM:2:ARG:HG3	12:AM:6:ILE:H	1.61	0.66
21:AN:58:ARG:H	21:AN:59:GLN:HE21	1.44	0.66
23:BB:2328:A:H2'	23:BB:2329:U:H6	1.58	0.66
50:BT:53:VAL:HG11	50:BT:87:LEU:HD13	1.78	0.66
1:CA:1004:A:H5'	1:CA:1024:G:N2	2.10	0.66
1:CA:1118:U:H1'	1:CA:1179:A:C5	2.31	0.66
1:CA:1206:G:H2'	1:CA:1207:G:O4'	1.95	0.66
2:CC:149:LYS:HG3	2:CC:168:ARG:HB2	1.77	0.66
36:D2:9:VAL:HG13	36:D2:10:LEU:H	1.60	0.66
23:DB:1119:U:H2'	23:DB:1120:G:H8	1.61	0.66
23:DB:1189:A:H2'	23:DB:1190:G:O4'	1.96	0.66
23:DB:2387:U:H1'	52:DW:38:ARG:CZ	2.26	0.66
23:DB:347:A:H2'	23:DB:348:A:H8	1.61	0.66
29:DE:58:LYS:HZ2	29:DE:58:LYS:H	1.42	0.66
47:DF:111:ARG:HE	47:DF:135:ILE:HG22	1.60	0.66
42:DN:47:VAL:C	42:DN:50:PRO:HD2	2.15	0.66
43:DO:100:HIS:O	43:DO:104:GLN:HB2	1.96	0.66
45:DS:84:ARG:HB3	45:DS:96:ILE:HG23	1.78	0.66
50:DT:5:GLU:O	50:DT:9:LYS:HE3	1.96	0.66
1:AA:518:C:H2'	1:AA:530:G:C8	2.31	0.66
4:AE:156:ARG:HD2	7:AH:42:GLU:O	1.95	0.66
23:BB:70:G:H3'	23:BB:113:U:H4'	1.76	0.66
23:BB:1437:C:H2'	23:BB:1438:U:H6	1.60	0.66
23:BB:17:G:H2'	23:BB:18:U:C6	2.30	0.66
23:BB:2537:U:H2'	23:BB:2538:C:C6	2.31	0.66
23:BB:974:G:H1'	23:BB:975:A:C8	2.31	0.66
29:BE:58:LYS:HB2	29:BE:60:TRP:CD1	2.31	0.66
1:CA:429:U:H1'	1:CA:430:A:H5''	1.77	0.66
1:CA:202:G:H21	1:CA:465:A:H61	1.44	0.66
1:CA:674:G:H2'	1:CA:675:A:C8	2.30	0.66
1:CA:923:A:H2'	1:CA:924:C:C6	2.31	0.66
3:CD:29:THR:H	3:CD:33:ILE:CG2	2.09	0.66
5:CF:1:MET:SD	5:CF:67:PRO:HD3	2.35	0.66
1:CA:1225:A:H5'	12:CM:101:THR:OG1	1.95	0.66
23:DB:1173:U:H1'	23:DB:1177:G:N1	2.11	0.66
23:DB:2010:G:H5''	45:DS:42:LYS:HB2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2064:C:H2'	23:DB:2065:C:C6	2.31	0.66
26:DD:8:LYS:HB2	26:DD:201:LEU:HD11	1.77	0.66
39:DX:13:GLU:HB2	39:DX:57:LEU:HD23	1.78	0.66
1:AA:1226:C:N4	12:AM:102:LYS:HB3	2.11	0.66
21:AN:31:SER:N	21:AN:44:VAL:HG11	2.11	0.66
23:BB:1138:G:H2'	23:BB:1139:G:O4'	1.95	0.66
23:BB:2841:C:H2'	23:BB:2842:G:H8	1.61	0.66
23:BB:979:A:H2'	23:BB:982:C:N4	2.10	0.66
40:BH:82:SER:N	40:BH:146:VAL:HG13	2.10	0.66
27:BK:20:MET:O	27:BK:41:ILE:HG13	1.95	0.66
51:BZ:54:LYS:HA	51:BZ:57:ARG:HD3	1.78	0.66
18:CB:107:ARG:HA	18:CB:110:ILE:HD12	1.78	0.66
6:CG:12:LEU:HD22	6:CG:13:PRO:HD2	1.78	0.66
9:CJ:86:ALA:HA	9:CJ:89:ARG:HD2	1.76	0.66
23:DB:1683:U:H2'	23:DB:1684:G:C8	2.31	0.66
23:DB:584:C:OP1	44:DQ:5:ARG:HB3	1.96	0.66
23:DB:71:A:H4'	23:DB:72:U:H5'	1.78	0.66
23:DB:910:A:H2'	23:DB:911:A:H8	1.60	0.66
1:AA:1252:A:H2'	1:AA:1253:G:O4'	1.95	0.65
1:AA:950:U:H2'	1:AA:951:G:C8	2.31	0.65
2:AC:99:GLN:O	2:AC:100:ILE:HB	1.95	0.65
10:AK:17:ASP:HA	10:AK:80:ASN:O	1.95	0.65
13:AP:67:ILE:HG12	13:AP:72:ALA:HB2	1.78	0.65
32:B4:3:VAL:HG23	32:B4:4:ARG:H	1.60	0.65
23:BB:1709:U:H2'	23:BB:1710:G:H8	1.60	0.65
23:BB:2366:A:H2'	23:BB:2367:G:O4'	1.96	0.65
23:BB:466:A:N3	23:BB:683:U:H1'	2.11	0.65
26:BD:33:ARG:NH1	26:BD:76:GLY:HA3	2.11	0.65
29:BE:105:LEU:HD11	29:BE:177:PRO:HG3	1.77	0.65
48:BG:18:ILE:HA	48:BG:23:ILE:HG13	1.76	0.65
40:BH:90:LEU:HB2	40:BH:123:ARG:HB3	1.77	0.65
46:BU:26:ASN:ND2	46:BU:34:ILE:HD12	2.11	0.65
35:BV:9:ARG:NH2	35:BV:12:GLN:HA	2.10	0.65
1:CA:736:C:H2'	1:CA:737:C:C6	2.31	0.65
23:DB:2824:C:H3'	23:DB:2825:G:H21	1.61	0.65
23:DB:358:U:H2'	23:DB:359:G:H8	1.61	0.65
23:DB:742:A:H2'	23:DB:743:A:H8	1.62	0.65
26:DD:33:ARG:NH1	26:DD:76:GLY:HA3	2.11	0.65
1:CA:1432:G:H5'	28:DP:105:LYS:HG2	1.78	0.65
44:DQ:111:LYS:HB2	49:DR:48:LYS:HZ2	1.60	0.65
35:DV:9:ARG:NH2	35:DV:12:GLN:HA	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:42:LEU:HD23	18:AB:45:THR:HB	1.77	0.65
3:AD:25:ARG:NH1	3:AD:26:ALA:HB2	2.11	0.65
3:AD:96:ARG:NH1	3:AD:133:SER:HA	2.10	0.65
5:AF:29:ILE:HG21	5:AF:64:VAL:HG11	1.78	0.65
21:AN:60:ARG:HG2	21:AN:62:ARG:HE	1.61	0.65
20:AO:63:ARG:O	20:AO:67:LEU:HG	1.95	0.65
19:AU:40:PRO:HA	19:AU:44:ARG:HD2	1.78	0.65
23:BB:532:A:H4'	23:BB:533:G:C8	2.31	0.65
47:BF:1:ALA:HB1	47:BF:4:HIS:HB3	1.78	0.65
48:BG:91:VAL:HG23	48:BG:92:GLY:H	1.62	0.65
27:BK:35:VAL:HG23	27:BK:36:GLY:N	2.07	0.65
1:CA:713:G:H2'	1:CA:714:G:C8	2.32	0.65
5:CF:52:ASN:HA	5:CF:53:LYS:HZ3	1.60	0.65
23:DB:1099:G:C5'	24:DI:3:LYS:N	2.59	0.65
23:DB:1704:C:H2'	23:DB:1705:A:C8	2.31	0.65
23:DB:320:A:H4'	23:DB:322:A:N7	2.11	0.65
25:DC:123:ILE:HD12	25:DC:191:LEU:HD11	1.77	0.65
25:DC:233:GLY:H	25:DC:241:LYS:NZ	1.95	0.65
48:DG:155:PRO:HA	48:DG:170:THR:HG22	1.77	0.65
40:DH:26:ALA:HA	40:DH:30:LEU:HB2	1.77	0.65
1:AA:1063:C:H3'	1:AA:1064:G:H2'	1.78	0.65
1:AA:193:C:H2'	1:AA:194:C:C6	2.31	0.65
1:AA:677:U:H2'	1:AA:678:U:C6	2.31	0.65
2:AC:155:ARG:H	2:AC:162:ALA:HA	1.61	0.65
21:AN:26:LEU:HA	21:AN:29:ILE:HD12	1.77	0.65
23:BB:171:U:H2'	23:BB:172:A:C8	2.31	0.65
23:BB:2306:C:H3'	23:BB:2307:G:H5'	1.78	0.65
47:BF:121:PHE:HA	47:BF:127:TYR:HA	1.77	0.65
22:BA:54:G:H21	47:BF:25:MET:HG2	1.60	0.65
49:BR:23:GLU:O	49:BR:25:LEU:HD22	1.97	0.65
39:BX:23:ARG:O	39:BX:27:ASN:HB2	1.97	0.65
1:CA:1006:G:H2'	1:CA:1007:U:H6	1.62	0.65
1:CA:1127:G:H2'	1:CA:1128:C:H6	1.61	0.65
1:CA:469:C:H2'	1:CA:470:C:H6	1.61	0.65
18:CB:16:GLY:HA2	18:CB:40:ILE:CD1	2.25	0.65
8:CI:56:MET:SD	8:CI:57:VAL:N	2.70	0.65
8:CI:87:MET:HB2	8:CI:94:ARG:CZ	2.27	0.65
23:DB:146:A:H2'	23:DB:147:C:C6	2.31	0.65
23:DB:1845:G:O2'	23:DB:1846:G:H5'	1.96	0.65
26:DD:107:VAL:HG13	26:DD:203:VAL:HG23	1.79	0.65
47:DF:141:ASP:O	47:DF:145:VAL:HG13	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DL:95:LEU:HB2	37:DL:101:ILE:CG1	2.26	0.65
44:DQ:89:ILE:HG22	44:DQ:91:ARG:H	1.60	0.65
12:AM:28:ARG:HH21	12:AM:62:PHE:HB2	1.60	0.65
16:AS:15:LEU:O	16:AS:18:VAL:HG12	1.97	0.65
23:BB:146:A:H2'	23:BB:147:C:C6	2.32	0.65
52:BW:39:GLN:NE2	52:BW:43:LYS:HB2	2.12	0.65
1:CA:1359:C:H3'	21:CN:74:ARG:NH2	2.11	0.65
1:CA:834:U:H2'	1:CA:835:U:C6	2.31	0.65
10:CK:124:LYS:HA	19:CU:34:ARG:HB3	1.77	0.65
12:CM:63:VAL:HG12	12:CM:68:LEU:HG	1.78	0.65
20:CO:11:ILE:HD11	20:CO:31:LEU:HD23	1.78	0.65
16:CS:31:ARG:HD2	16:CS:31:ARG:H	1.61	0.65
17:CT:38:ILE:HD11	17:CT:82:ILE:HG22	1.77	0.65
19:CU:38:GLU:C	19:CU:40:PRO:HD2	2.16	0.65
33:D1:8:ILE:CD1	33:D1:51:ALA:HA	2.25	0.65
48:DG:36:LEU:H	48:DG:36:LEU:HD22	1.61	0.65
27:DK:19:VAL:C	27:DK:41:ILE:HD11	2.16	0.65
1:AA:1063:C:N4	1:AA:1193:G:H1	1.93	0.65
1:AA:812:G:O2'	1:AA:813:U:H6	1.78	0.65
2:AC:111:ASP:HB3	2:AC:114:LEU:HD12	1.77	0.65
2:AC:77:GLY:HA3	2:AC:81:GLU:HB3	1.79	0.65
5:AF:72:ASP:HA	5:AF:75:GLU:OE1	1.95	0.65
6:AG:100:MET:HA	6:AG:103:ILE:HD12	1.76	0.65
8:AI:12:LYS:N	8:AI:105:ARG:HH21	1.93	0.65
23:BB:1461:C:H2'	23:BB:1462:C:H6	1.62	0.65
23:BB:1480:C:H2'	23:BB:1481:U:O4'	1.95	0.65
23:BB:181:A:H2'	23:BB:182:A:C8	2.32	0.65
23:BB:2591:C:H2'	23:BB:2592:G:H8	1.62	0.65
26:BD:107:VAL:HG13	26:BD:203:VAL:HG23	1.76	0.65
26:BD:55:LYS:HB3	26:BD:55:LYS:HZ2	1.61	0.65
29:BE:46:GLN:HB3	29:BE:86:ALA:HA	1.79	0.65
48:BG:166:GLU:HG2	48:BG:168:VAL:HG23	1.77	0.65
28:BP:89:GLY:HA2	28:BP:112:ARG:N	2.12	0.65
46:BU:40:LEU:H	46:BU:40:LEU:HD12	1.61	0.65
1:CA:1298:U:H4'	1:CA:1299:A:O4'	1.95	0.65
1:CA:1333:A:H2'	1:CA:1334:G:O4'	1.97	0.65
1:CA:715:A:H2'	1:CA:716:A:H8	1.60	0.65
1:CA:620:C:C2	3:CD:131:ILE:HD13	2.31	0.65
4:CE:113:VAL:HG23	4:CE:114:LEU:N	2.10	0.65
5:CF:6:ILE:HG23	5:CF:62:MET:HB3	1.78	0.65
10:CK:58:THR:HB	10:CK:59:PRO:HD2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:CN:20:PHE:CA	21:CN:24:ALA:HB2	2.27	0.65
23:DB:1173:U:H1'	23:DB:1177:G:H1	1.62	0.65
23:DB:171:U:H2'	23:DB:172:A:C8	2.32	0.65
23:DB:466:A:N3	23:DB:683:U:H1'	2.11	0.65
23:DB:851:C:O4'	30:DY:46:MET:HG2	1.97	0.65
35:DV:72:VAL:HB	35:DV:92:VAL:O	1.96	0.65
1:AA:1043:G:H2'	1:AA:1044:A:C8	2.32	0.65
1:AA:80:A:C4	1:AA:81:A:H1'	2.32	0.65
4:AE:131:ASN:HD22	4:AE:134:ASN:H	1.44	0.65
8:AI:25:GLY:HA2	8:AI:60:LEU:O	1.96	0.65
11:AL:64:SER:OG	11:AL:96:THR:HG23	1.97	0.65
23:BB:1551:A:H3'	23:BB:1552:A:H5''	1.78	0.65
23:BB:1724:G:H2'	23:BB:1725:U:C6	2.31	0.65
23:BB:181:A:H2'	23:BB:182:A:H8	1.60	0.65
23:BB:228:C:H4'	23:BB:229:C:H5''	1.78	0.65
23:BB:565:C:H2'	23:BB:566:U:C6	2.32	0.65
23:BB:910:A:H2'	23:BB:911:A:C8	2.32	0.65
48:BG:30:GLY:CA	48:BG:78:VAL:HA	2.27	0.65
27:BK:38:ILE:N	27:BK:38:ILE:HD13	2.12	0.65
38:BM:10:ARG:HH11	38:BM:89:VAL:HG22	1.62	0.65
51:BZ:56:MET:HA	51:BZ:59:ILE:HG12	1.78	0.65
21:CN:20:PHE:HA	21:CN:24:ALA:HB2	1.78	0.65
13:CP:54:LEU:HD13	13:CP:80:LYS:HE3	1.78	0.65
23:DB:328:U:H4'	46:DU:65:GLN:NE2	2.10	0.65
23:DB:565:C:H2'	23:DB:566:U:C6	2.31	0.65
23:DB:784:G:O2'	23:DB:785:G:H5''	1.96	0.65
23:DB:845:A:C2	23:DB:847:U:H1'	2.31	0.65
25:DC:93:VAL:O	25:DC:94:LEU:HB3	1.97	0.65
48:DG:166:GLU:H	48:DG:166:GLU:CD	2.00	0.65
18:AB:46:VAL:CG1	18:AB:47:PRO:HD3	2.25	0.65
10:AK:58:THR:HB	10:AK:59:PRO:HD2	1.78	0.65
22:BA:91:C:H2'	22:BA:92:C:C6	2.32	0.65
23:BB:1139:G:O2'	23:BB:1140:C:H5'	1.96	0.65
23:BB:1440:U:H2'	23:BB:1441:G:H8	1.62	0.65
23:BB:2776:A:H4'	23:BB:2777:G:C5'	2.26	0.65
48:BG:166:GLU:H	48:BG:166:GLU:CD	1.99	0.65
23:DB:2678:C:H2'	23:DB:2679:A:H8	1.60	0.65
23:DB:627:A:H4'	23:DB:628:G:H5'	1.78	0.65
23:DB:942:G:H2'	23:DB:943:A:O4'	1.96	0.65
41:DJ:11:VAL:HG11	41:DJ:13:ARG:HE	1.60	0.65
52:DW:76:ARG:NH2	52:DW:76:ARG:HA	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:543:U:H2'	1:AA:544:G:H8	1.62	0.65
18:AB:163:ILE:HG23	18:AB:164:ASP:N	2.11	0.65
18:AB:76:SER:HA	18:AB:92:ASN:ND2	2.12	0.65
2:AC:52:SER:HA	2:AC:113:LYS:HG2	1.79	0.65
10:AK:52:ARG:HB3	10:AK:52:ARG:HH11	1.61	0.65
15:AR:60:ARG:HA	15:AR:63:TYR:HD2	1.62	0.65
17:AT:43:LYS:HB3	17:AT:86:ALA:HB1	1.79	0.65
23:BB:1309:G:OP1	36:B2:9:VAL:HG12	1.97	0.65
23:BB:1429:G:H2'	23:BB:1430:G:H8	1.62	0.65
23:BB:165:A:H2'	23:BB:166:U:C6	2.32	0.65
23:BB:2230:G:H2'	23:BB:2231:U:C6	2.31	0.65
41:BJ:43:GLU:O	41:BJ:45:THR:N	2.30	0.65
37:BL:79:LEU:HB2	37:BL:113:ALA:H	1.60	0.65
51:BZ:77:LYS:HD3	51:BZ:77:LYS:H	1.60	0.65
1:CA:880:C:H2'	1:CA:881:G:H8	1.60	0.65
18:CB:151:LYS:HG3	18:CB:152:ASP:N	2.11	0.65
2:CC:41:TYR:HD1	2:CC:42:LEU:HD22	1.62	0.65
10:CK:28:ASN:ND2	10:CK:29:THR:H	1.95	0.65
36:D2:13:ASN:O	36:D2:17:GLY:N	2.30	0.65
22:DA:91:C:H2'	22:DA:92:C:H6	1.62	0.65
47:DF:163:GLU:HA	47:DF:166:ARG:HD2	1.77	0.65
48:DG:154:GLU:HG2	48:DG:156:TYR:H	1.62	0.65
40:DH:4:ILE:HB	40:DH:37:VAL:CG1	2.26	0.65
35:DV:72:VAL:HG12	35:DV:93:ARG:HA	1.79	0.65
18:AB:49:PHE:HA	18:AB:212:TYR:OH	1.97	0.65
18:AB:224:ARG:O	18:AB:225:SER:HB2	1.96	0.65
18:AB:61:SER:HB2	18:AB:62:ARG:HH11	1.62	0.65
23:BB:2469:A:H4'	38:BM:55:ARG:NE	2.12	0.65
23:BB:2615:U:H1'	31:B0:3:GLN:HB3	1.79	0.65
26:BD:159:LYS:HZ2	26:BD:160:LYS:N	1.95	0.65
26:BD:40:LEU:HD23	26:BD:46:ARG:HG2	1.78	0.65
47:BF:141:ASP:HB3	47:BF:144:LYS:HB2	1.79	0.65
47:BF:2:LYS:CE	47:BF:2:LYS:H	2.10	0.65
23:BB:1199:U:H5'	44:BQ:4:LYS:HD3	1.77	0.65
30:BY:8:GLN:HG2	30:BY:31:ILE:HA	1.78	0.65
1:CA:1361:G:H2'	1:CA:1362:A:H5''	1.77	0.65
2:CC:149:LYS:HB2	2:CC:168:ARG:HG3	1.79	0.65
6:CG:90:VAL:HB	6:CG:94:ARG:HD2	1.79	0.65
22:DA:42:C:H4'	47:DF:63:LYS:HB3	1.79	0.65
23:DB:1534:U:H2'	23:DB:1536:C:N3	2.12	0.65
23:DB:215:G:H4'	23:DB:216:A:H4'	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2471:A:O2'	23:DB:2472:G:H8	1.80	0.65
25:DC:244:VAL:HB	25:DC:249:VAL:N	2.06	0.65
29:DE:146:VAL:HA	29:DE:185:LYS:O	1.96	0.65
27:DK:35:VAL:HG23	27:DK:36:GLY:N	2.08	0.65
37:DL:129:LYS:HA	37:DL:132:ARG:HG2	1.79	0.65
43:DO:59:ALA:HA	43:DO:62:LEU:HG	1.77	0.65
23:BB:1821:A:H2'	23:BB:1822:C:C6	2.32	0.65
23:BB:2071:A:H2'	23:BB:2072:C:C6	2.32	0.65
23:BB:2196:C:O2'	23:BB:2197:U:H5'	1.97	0.65
23:BB:2720:U:H5''	28:BP:52:ARG:NH2	2.12	0.65
48:BG:153:PRO:HB3	48:BG:158:GLY:HA2	1.77	0.65
39:BX:48:ARG:O	39:BX:51:ALA:HB3	1.97	0.65
51:BZ:40:VAL:HG22	51:BZ:45:ARG:O	1.97	0.65
6:CG:14:ASP:CG	6:CG:17:PHE:HB2	2.17	0.65
7:CH:86:LYS:HD2	7:CH:90:GLU:HG2	1.79	0.65
36:D2:27:GLY:O	36:D2:30:VAL:HB	1.97	0.65
23:DB:2861:U:H2'	23:DB:2862:G:H8	1.61	0.65
23:DB:362:A:H3'	23:DB:363:G:H8	1.61	0.65
29:DE:67:ARG:HG2	29:DE:67:ARG:HH11	1.62	0.65
29:DE:46:GLN:HB3	29:DE:86:ALA:HA	1.79	0.65
48:DG:166:GLU:HG2	48:DG:168:VAL:HG23	1.77	0.65
28:DP:20:ARG:HG3	28:DP:21:PRO:CD	2.27	0.65
30:DY:9:THR:HB	30:DY:53:MET:O	1.96	0.65
1:AA:620:C:C2	3:AD:131:ILE:HD13	2.32	0.64
11:AL:49:ARG:HB2	11:AL:89:LEU:HD11	1.79	0.64
23:BB:1007:C:H4'	41:BJ:110:PRO:HB3	1.79	0.64
23:BB:2142:A:O2'	23:BB:2143:C:H5'	1.97	0.64
23:BB:2824:C:H3'	23:BB:2825:G:N2	2.13	0.64
23:BB:448:U:C5	23:BB:583:G:H1'	2.32	0.64
48:BG:21:GLN:OE1	48:BG:37:ASN:HB2	1.96	0.64
40:BH:114:GLU:HB2	40:BH:132:PHE:CD1	2.32	0.64
40:BH:26:ALA:HA	40:BH:30:LEU:HB2	1.78	0.64
27:BK:19:VAL:HG12	27:BK:41:ILE:HG12	1.79	0.64
35:BV:72:VAL:HG12	35:BV:93:ARG:HA	1.80	0.64
6:CG:75:LYS:HA	6:CG:75:LYS:NZ	2.12	0.64
12:CM:12:LYS:HD2	12:CM:16:ILE:HD13	1.78	0.64
23:DB:1021:A:H62	23:DB:1141:U:H3	1.45	0.64
23:DB:1857:G:N2	23:DB:1884:G:H2'	2.12	0.64
23:DB:2776:A:H4'	23:DB:2777:G:C5'	2.26	0.64
23:DB:351:C:H2'	23:DB:352:A:C8	2.31	0.64
29:DE:132:LYS:O	29:DE:135:ALA:HB3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DK:60:ALA:HB2	27:DK:86:LEU:HA	1.77	0.64
37:DL:111:ILE:HG22	37:DL:112:LEU:H	1.62	0.64
38:DM:105:MET:HB2	38:DM:117:PHE:HZ	1.61	0.64
52:DW:23:LYS:HD2	52:DW:24:ARG:H	1.61	0.64
30:DY:8:GLN:HG2	30:DY:31:ILE:HA	1.77	0.64
1:AA:803:G:H2'	1:AA:804:U:C6	2.33	0.64
9:AJ:52:LEU:HB2	21:AN:80:ARG:HE	1.60	0.64
22:BA:25:U:H4'	22:BA:26:C:H5''	1.77	0.64
23:BB:570:G:H2'	23:BB:2030:A:N7	2.11	0.64
23:BB:2752:C:H2'	23:BB:2753:A:O4'	1.97	0.64
23:BB:654:A:N3	23:BB:655:A:H5''	2.12	0.64
40:BH:89:LYS:HA	40:BH:89:LYS:NZ	2.11	0.64
24:BI:129:GLU:HB3	24:BI:133:ARG:NH1	2.10	0.64
49:BR:39:LEU:HB3	49:BR:49:ILE:HD11	1.78	0.64
50:BT:50:LEU:HD22	50:BT:50:LEU:H	1.61	0.64
52:BW:50:VAL:HG23	52:BW:61:LYS:HE3	1.78	0.64
30:BY:23:LEU:HD13	30:BY:28:LEU:HB2	1.78	0.64
18:CB:61:SER:HB2	18:CB:62:ARG:NH1	2.12	0.64
18:CB:65:LYS:HB2	18:CB:158:ASP:CG	2.18	0.64
8:CI:11:ARG:HA	8:CI:105:ARG:CZ	2.27	0.64
11:CL:14:LYS:HZ3	11:CL:16:ALA:HB3	1.62	0.64
19:CU:36:PHE:HB3	19:CU:40:PRO:CD	2.27	0.64
23:DB:1440:U:H2'	23:DB:1441:G:H8	1.62	0.64
23:DB:1461:C:H2'	23:DB:1462:C:H6	1.62	0.64
23:DB:17:G:H2'	23:DB:18:U:C6	2.31	0.64
23:DB:2230:G:H2'	23:DB:2231:U:C6	2.31	0.64
23:DB:932:U:H5'	23:DB:933:A:OP1	1.97	0.64
26:DD:111:GLY:H	26:DD:194:PRO:HG2	1.62	0.64
1:AA:564:C:C2	14:AQ:32:ILE:HD11	2.32	0.64
5:AF:52:ASN:HA	5:AF:53:LYS:HZ3	1.63	0.64
23:BB:1021:A:H62	23:BB:1141:U:H3	1.43	0.64
23:BB:1857:G:N2	23:BB:1884:G:H2'	2.12	0.64
23:BB:2841:C:H2'	23:BB:2842:G:C8	2.31	0.64
23:BB:63:A:H2'	23:BB:64:A:N7	2.12	0.64
38:BM:12:MET:HB2	38:BM:72:PRO:HG2	1.78	0.64
1:CA:1041:G:H2'	1:CA:1042:A:C8	2.31	0.64
1:CA:194:C:O2'	1:CA:195:A:H5'	1.96	0.64
16:CS:51:HIS:HB2	16:CS:56:HIS:CD2	2.31	0.64
23:DB:2216:G:H2'	23:DB:2217:G:H8	1.62	0.64
23:DB:2469:A:H2'	23:DB:2470:G:O4'	1.98	0.64
26:DD:33:ARG:HE	26:DD:51:THR:HB	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:54:ALA:HA	26:DD:76:GLY:HA2	1.78	0.64
40:DH:114:GLU:HG3	40:DH:133:GLN:O	1.96	0.64
27:DK:20:MET:O	27:DK:41:ILE:HG13	1.98	0.64
37:DL:121:THR:HB	37:DL:141:LYS:HB3	1.79	0.64
44:DQ:105:PHE:HA	44:DQ:108:LEU:HB2	1.78	0.64
1:AA:1134:G:C2	1:AA:1135:U:H1'	2.33	0.64
1:AA:1226:C:H4'	1:AA:1227:A:OP1	1.98	0.64
1:AA:1461:G:H2'	1:AA:1462:C:C6	2.32	0.64
1:AA:513:C:H2'	1:AA:514:C:H6	1.62	0.64
18:AB:112:ARG:HE	18:AB:116:LEU:HD11	1.63	0.64
18:AB:187:ASP:CG	18:AB:203:ASP:HB3	2.18	0.64
33:B1:32:LYS:HA	33:B1:51:ALA:O	1.97	0.64
32:B4:16:ILE:HA	32:B4:25:VAL:HG22	1.79	0.64
22:BA:42:C:H4'	47:BF:63:LYS:HB3	1.80	0.64
23:BB:155:A:H2'	23:BB:156:A:C8	2.33	0.64
23:BB:2892:G:H5''	23:BB:2894:G:N2	2.13	0.64
23:BB:634:C:H2'	23:BB:635:C:H6	1.61	0.64
24:BI:25:PRO:O	24:BI:29:GLN:HG2	1.98	0.64
38:BM:33:LEU:HD22	38:BM:128:THR:HB	1.78	0.64
1:CA:382:A:H2'	1:CA:383:A:C8	2.32	0.64
1:CA:940:C:H2'	1:CA:941:G:H8	1.62	0.64
3:CD:172:VAL:HA	3:CD:179:GLY:HA2	1.78	0.64
8:CI:11:ARG:NH2	8:CI:12:LYS:HB2	2.12	0.64
11:CL:13:ARG:O	11:CL:14:LYS:HB3	1.96	0.64
15:CR:31:TYR:O	15:CR:39:VAL:HG22	1.97	0.64
23:DB:1485:U:H2'	23:DB:1486:U:C6	2.33	0.64
23:DB:1843:C:H5''	25:DC:250:GLN:HE21	1.61	0.64
23:DB:558:U:OP1	41:DJ:114:LEU:HB2	1.97	0.64
48:DG:89:VAL:HG12	48:DG:90:GLY:H	1.60	0.64
40:DH:116:ARG:HB3	40:DH:131:SER:N	2.13	0.64
23:DB:1132:U:O2	41:DJ:75:TYR:HB2	1.98	0.64
38:DM:41:LEU:HB3	38:DM:46:ILE:CG2	2.27	0.64
35:DV:14:LYS:HE3	35:DV:18:ARG:HH21	1.62	0.64
1:AA:17:U:H2'	1:AA:18:C:H6	1.62	0.64
1:AA:250:A:H1'	1:AA:252:U:C6	2.32	0.64
2:AC:19:SER:HB3	2:AC:21:TRP:HE1	1.62	0.64
3:AD:2:ARG:NH1	3:AD:114:ARG:HD3	2.12	0.64
6:AG:71:THR:H	6:AG:141:HIS:CE1	2.14	0.64
6:AG:149:ALA:H	10:AK:55:ARG:HH22	1.43	0.64
22:BA:87:U:H2'	22:BA:88:C:O5'	1.97	0.64
37:BL:129:LYS:HA	37:BL:132:ARG:HG2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:7:LEU:C	50:BT:9:LYS:HD2	2.18	0.64
46:BU:78:LYS:HD3	46:BU:79:ALA:H	1.63	0.64
18:CB:59:ILE:HG22	18:CB:62:ARG:NE	2.12	0.64
2:CC:151:GLU:HB3	2:CC:198:LYS:HB2	1.79	0.64
1:CA:408:A:OP1	3:CD:111:ALA:HB3	1.97	0.64
6:CG:50:ALA:HB2	6:CG:57:GLU:HG3	1.80	0.64
4:CE:156:ARG:HD2	7:CH:42:GLU:O	1.98	0.64
21:CN:60:ARG:HD3	21:CN:62:ARG:NH2	2.12	0.64
1:CA:564:C:C2	14:CQ:32:ILE:HD11	2.33	0.64
23:DB:1406:U:H2'	23:DB:1407:G:H8	1.62	0.64
23:DB:2071:A:H2'	23:DB:2072:C:C6	2.33	0.64
23:DB:289:G:H2'	23:DB:290:U:C6	2.31	0.64
23:DB:2336:A:H61	52:DW:40:ARG:HG3	1.60	0.64
18:AB:95:TRP:HZ2	18:AB:100:LEU:HD13	1.63	0.64
23:BB:2698:U:H2'	23:BB:2699:C:C6	2.32	0.64
23:BB:287:G:H2'	23:BB:288:U:C6	2.33	0.64
23:BB:324:A:H2'	23:BB:325:G:O4'	1.97	0.64
48:BG:154:GLU:HG2	48:BG:156:TYR:H	1.62	0.64
23:BB:2356:U:H5'	52:BW:16:GLU:HG3	1.79	0.64
23:DB:1098:A:OP2	24:DI:3:LYS:HG2	1.98	0.64
23:DB:1149:G:H2'	23:DB:1150:C:C6	2.32	0.64
23:DB:1778:U:H2'	23:DB:1784:A:H62	1.63	0.64
23:DB:2008:C:H2'	23:DB:2009:A:C8	2.30	0.64
26:DD:10:GLY:HA3	26:DD:26:VAL:N	2.06	0.64
47:DF:169:LEU:HA	47:DF:172:PHE:HD2	1.63	0.64
1:AA:1215:G:H5'	1:AA:1215:G:H8	1.61	0.64
1:AA:840:C:C5	1:AA:842:U:H4'	2.32	0.64
1:AA:91:U:H2'	1:AA:92:U:C6	2.32	0.64
2:AC:139:ASN:O	2:AC:143:LEU:HD23	1.97	0.64
10:AK:28:ASN:ND2	10:AK:29:THR:H	1.94	0.64
23:BB:1683:U:H2'	23:BB:1684:G:H8	1.61	0.64
23:BB:2241:A:H2'	23:BB:2242:G:C8	2.33	0.64
23:BB:322:A:C5'	23:BB:340:A:H1'	2.19	0.64
23:BB:946:C:H2'	23:BB:947:A:H8	1.62	0.64
27:BK:60:ALA:HB2	27:BK:86:LEU:HA	1.77	0.64
43:BO:67:ASN:HB3	43:BO:70:ALA:HB2	1.79	0.64
44:BQ:111:LYS:HB2	49:BR:48:LYS:HZ2	1.63	0.64
1:CA:1317:C:H4'	21:CN:48:GLN:OE1	1.97	0.64
1:CA:522:C:H41	11:CL:49:ARG:HH22	1.46	0.64
1:CA:598:U:H2'	1:CA:599:C:C6	2.33	0.64
15:CR:72:ARG:HH11	15:CR:72:ARG:N	1.84	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2306:C:H3'	23:DB:2307:G:H5'	1.79	0.64
23:DB:2678:C:H2'	23:DB:2679:A:C8	2.32	0.64
23:DB:303:G:H2'	23:DB:304:U:C6	2.33	0.64
23:DB:718:A:H3'	23:DB:719:C:C6	2.33	0.64
23:DB:876:C:C4	23:DB:877:A:H1'	2.32	0.64
23:DB:2575:C:H5''	26:DD:149:ASN:HD22	1.63	0.64
47:DF:141:ASP:HB3	47:DF:144:LYS:HB2	1.79	0.64
48:DG:144:ALA:HB1	48:DG:163:TYR:HE1	1.62	0.64
49:DR:39:LEU:CB	49:DR:49:ILE:HD11	2.28	0.64
50:DT:69:ARG:HG2	50:DT:73:ARG:C	2.18	0.64
52:DW:50:VAL:HG23	52:DW:61:LYS:HE3	1.80	0.64
1:AA:1314:C:N4	16:AS:3:SER:HB3	2.12	0.64
1:AA:408:A:OP1	3:AD:111:ALA:HB3	1.98	0.64
1:AA:642:A:H2'	1:AA:643:C:H6	1.63	0.64
18:AB:120:SER:HA	18:AB:125:PHE:HB2	1.77	0.64
6:AG:74:VAL:HG21	6:AG:85:GLN:HE21	1.62	0.64
8:AI:20:ILE:HG12	8:AI:86:LEU:HD12	1.80	0.64
10:AK:88:PRO:HA	10:AK:92:ARG:HD2	1.80	0.64
11:AL:31:GLY:HA3	11:AL:54:VAL:CG1	2.28	0.64
21:AN:27:LYS:HG3	21:AN:28:ALA:H	1.61	0.64
19:AU:38:GLU:C	19:AU:40:PRO:HD2	2.18	0.64
23:BB:2099:U:H2'	23:BB:2100:G:H8	1.63	0.64
23:BB:620:G:H5'	23:BB:620:G:N3	2.13	0.64
23:BB:71:A:H4'	23:BB:72:U:H5'	1.79	0.64
41:BJ:93:ILE:HA	41:BJ:97:PRO:HG3	1.79	0.64
1:CA:1023:U:H2'	1:CA:1024:G:O4'	1.97	0.64
1:CA:335:C:H2'	1:CA:336:A:H8	1.63	0.64
1:CA:642:A:H2'	1:CA:643:C:H6	1.63	0.64
18:CB:159:ALA:HB1	18:CB:183:PHE:CE1	2.26	0.64
18:CB:27:LYS:HB3	18:CB:28:PRO:HD3	1.80	0.64
32:D4:5:ALA:HA	32:D4:37:GLN:NE2	2.12	0.64
23:DB:1429:G:H2'	23:DB:1430:G:H8	1.63	0.64
23:DB:2794:C:H2'	23:DB:2795:C:C6	2.33	0.64
23:DB:533:G:H2'	23:DB:534:U:C6	2.33	0.64
23:DB:947:A:H2'	23:DB:948:C:H6	1.62	0.64
25:DC:143:VAL:HB	25:DC:153:LEU:HB2	1.80	0.64
25:DC:233:GLY:H	25:DC:241:LYS:HZ2	1.42	0.64
47:DF:32:LYS:HA	47:DF:95:MET:HG3	1.80	0.64
41:DJ:93:ILE:HA	41:DJ:97:PRO:HG3	1.79	0.64
50:DT:50:LEU:HD22	50:DT:50:LEU:H	1.62	0.64
1:AA:1147:C:H2'	1:AA:1148:U:H6	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:157:U:O2'	1:AA:158:G:H5'	1.98	0.64
1:AA:713:G:H2'	1:AA:714:G:C8	2.33	0.64
8:AI:14:SER:HA	8:AI:68:GLY:C	2.17	0.64
20:AO:85:LEU:HB3	20:AO:87:LEU:HG	1.80	0.64
23:BB:2773:C:H5''	26:BD:169:ARG:HB2	1.79	0.64
23:BB:2898:U:H2'	23:BB:2899:A:C8	2.33	0.64
23:BB:857:G:O2'	23:BB:858:G:H5'	1.98	0.64
26:BD:118:PHE:O	26:BD:119:ALA:HB3	1.97	0.64
40:BH:100:ALA:CB	40:BH:112:LYS:HA	2.28	0.64
40:BH:94:ILE:HB	40:BH:128:HIS:NE2	2.13	0.64
43:BO:59:ALA:HA	43:BO:62:LEU:HG	1.79	0.64
28:BP:59:THR:H	28:BP:72:VAL:HA	1.62	0.64
45:BS:26:GLY:N	45:BS:71:VAL:HG13	2.11	0.64
1:CA:1032:G:H3'	1:CA:1032:G:N3	2.12	0.64
1:CA:1097:C:H2'	1:CA:1098:C:H6	1.63	0.64
1:CA:1512:U:H2'	1:CA:1513:A:H8	1.63	0.64
1:CA:18:C:H1'	1:CA:1079:G:H21	1.62	0.64
1:CA:840:C:C5	1:CA:842:U:H4'	2.33	0.64
23:DB:1724:G:H2'	23:DB:1725:U:C6	2.32	0.64
23:DB:184:C:H2'	23:DB:185:G:C8	2.32	0.64
47:DF:134:GLN:H	47:DF:150:GLY:N	1.94	0.64
27:DK:70:ARG:HD3	27:DK:76:VAL:HG22	1.80	0.64
28:DP:88:ARG:HB2	28:DP:112:ARG:NH1	2.12	0.64
50:DT:34:VAL:HG11	50:DT:43:ILE:HD11	1.79	0.64
1:AA:1234:C:O2'	1:AA:1235:U:H5'	1.98	0.64
1:AA:301:G:H2'	1:AA:302:G:C8	2.32	0.64
1:AA:704:A:C2	1:AA:705:G:H1'	2.33	0.64
1:AA:923:A:OP1	4:AE:25:LYS:HB3	1.98	0.64
16:AS:49:ALA:O	16:AS:56:HIS:HB3	1.98	0.64
17:AT:74:HIS:O	17:AT:78:LEU:HG	1.98	0.64
23:BB:1149:G:H2'	23:BB:1150:C:C6	2.33	0.64
23:BB:2469:A:H2'	23:BB:2470:G:O4'	1.98	0.64
23:BB:580:U:H2'	23:BB:581:C:H6	1.63	0.64
26:BD:62:LYS:HB2	26:BD:63:PRO:HD3	1.80	0.64
47:BF:137:PHE:O	47:BF:139:GLU:N	2.31	0.64
41:BJ:109:LEU:HD13	41:BJ:119:PHE:HB2	1.79	0.64
1:CA:986:U:H2'	1:CA:987:G:O4'	1.97	0.64
18:CB:102:ASN:HD21	18:CB:105:THR:HB	1.60	0.64
2:CC:29:ALA:HB2	21:CN:75:LYS:O	1.98	0.64
1:CA:1216:A:H5''	21:CN:4:SER:HB2	1.79	0.64
21:CN:65:GLN:HE22	21:CN:78:LEU:HG	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1993:U:H4'	26:DD:133:THR:CG2	2.28	0.64
23:DB:1:G:H2'	23:DB:2:G:C8	2.33	0.64
40:DH:5:LEU:HB3	40:DH:13:GLY:HA2	1.80	0.64
27:DK:110:GLU:HA	27:DK:113:MET:HG3	1.80	0.64
27:DK:87:LEU:HD12	27:DK:93:GLN:N	2.12	0.64
49:DR:23:GLU:O	49:DR:25:LEU:HD22	1.98	0.64
1:AA:1019:A:H2'	1:AA:1020:G:O4'	1.98	0.63
1:AA:1342:C:H2'	1:AA:1343:G:C8	2.32	0.63
7:AH:23:ALA:HA	7:AH:62:LEU:H	1.62	0.63
16:AS:62:THR:HG22	16:AS:63:ASP:H	1.61	0.63
33:B1:7:LYS:HD3	34:B3:33:THR:HG21	1.81	0.63
23:BB:1485:U:H2'	23:BB:1486:U:C6	2.33	0.63
23:BB:2358:A:H61	37:BL:54:GLN:HE22	1.47	0.63
23:BB:705:A:N6	23:BB:726:G:H1'	2.14	0.63
26:BD:5:VAL:N	26:BD:32:ASN:HD21	1.96	0.63
37:BL:57:LEU:HD12	37:BL:60:ARG:NH2	2.13	0.63
18:CB:218:ALA:HA	18:CB:221:ARG:NH1	2.12	0.63
18:CB:57:ASN:CG	18:CB:219:THR:HB	2.18	0.63
16:CS:45:GLY:HA2	16:CS:60:PHE:HB3	1.80	0.63
33:D1:31:GLU:H	33:D1:31:GLU:CD	2.01	0.63
23:DB:2267:A:H3'	23:DB:2267:A:H8	1.63	0.63
25:DC:16:VAL:HB	25:DC:203:VAL:HG11	1.81	0.63
25:DC:6:LYS:O	25:DC:8:THR:HG23	1.97	0.63
26:DD:118:PHE:O	26:DD:119:ALA:HB3	1.98	0.63
26:DD:2:ILE:HG22	26:DD:82:PHE:HB3	1.79	0.63
23:DB:1100:C:OP2	24:DI:2:LYS:HB3	1.98	0.63
41:DJ:4:PHE:HB3	41:DJ:44:TYR:CD1	2.33	0.63
41:DJ:56:VAL:HG12	41:DJ:57:LEU:H	1.63	0.63
1:AA:860:A:H2'	1:AA:861:G:O4'	1.98	0.63
6:AG:17:PHE:HE2	6:AG:58:LEU:HD22	1.64	0.63
10:AK:75:GLU:H	10:AK:75:GLU:CD	2.01	0.63
22:BA:49:C:H2'	22:BA:50:A:C8	2.32	0.63
23:BB:1024:G:C3'	23:BB:1025:G:H5''	2.27	0.63
23:BB:718:A:H3'	23:BB:719:C:C6	2.30	0.63
23:BB:764:A:H5''	25:BC:208:GLY:HA2	1.81	0.63
48:BG:104:LEU:HD22	48:BG:106:LEU:HD21	1.79	0.63
42:BN:33:ILE:HD11	42:BN:112:TYR:HD1	1.62	0.63
45:BS:17:VAL:C	45:BS:19:LEU:H	2.01	0.63
45:BS:66:ILE:N	45:BS:66:ILE:HD13	2.14	0.63
50:BT:54:GLU:HB3	50:BT:88:LYS:HB2	1.78	0.63
1:CA:1091:U:H3'	6:CG:3:ARG:NH1	2.12	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:394:G:H2'	1:CA:395:C:C6	2.32	0.63
1:CA:927:G:N2	1:CA:1391:U:H1'	2.13	0.63
18:CB:113:LEU:HD21	18:CB:151:LYS:HB2	1.78	0.63
12:CM:1:ALA:N	12:CM:8:ILE:HG22	2.13	0.63
23:DB:152:A:H2'	23:DB:153:U:C6	2.33	0.63
23:DB:1812:U:H2'	23:DB:1813:G:C8	2.33	0.63
25:DC:156:SER:O	25:DC:194:VAL:HG11	1.98	0.63
26:DD:10:GLY:O	26:DD:11:MET:HB2	1.98	0.63
26:DD:148:GLN:CB	26:DD:152:PRO:HG2	2.26	0.63
26:DD:4:LEU:HG	26:DD:29:VAL:HG11	1.80	0.63
47:DF:22:ASN:H	47:DF:22:ASN:ND2	1.96	0.63
27:DK:38:ILE:HD13	27:DK:38:ILE:N	2.13	0.63
50:DT:53:VAL:HG11	50:DT:87:LEU:HD13	1.79	0.63
1:AA:1324:A:H2'	1:AA:1325:C:H6	1.61	0.63
1:AA:98:A:H2'	1:AA:99:C:C6	2.34	0.63
3:AD:12:ARG:NH1	3:AD:12:ARG:HB3	2.12	0.63
5:AF:71:ILE:HG13	5:AF:72:ASP:N	2.13	0.63
23:BB:150:U:H2'	23:BB:151:C:C6	2.33	0.63
23:BB:1796:U:H2'	23:BB:1797:G:C8	2.34	0.63
23:BB:2471:A:O2'	23:BB:2472:G:H8	1.80	0.63
23:BB:2757:A:H2	48:BG:63:GLN:HE22	1.46	0.63
26:BD:119:ALA:HB1	26:BD:163:GLY:H	1.61	0.63
29:BE:33:VAL:O	29:BE:36:ALA:HB3	1.99	0.63
27:BK:25:LEU:HG	27:BK:39:ILE:HA	1.78	0.63
44:BQ:108:LEU:HA	49:BR:48:LYS:HD3	1.78	0.63
44:BQ:111:LYS:HB2	49:BR:48:LYS:NZ	2.12	0.63
1:CA:1036:A:H2'	1:CA:1037:C:C6	2.32	0.63
1:CA:858:G:O6	1:CA:869:G:H3'	1.99	0.63
23:DB:155:A:H2'	23:DB:156:A:C8	2.34	0.63
23:DB:1709:U:H2'	23:DB:1710:G:C8	2.33	0.63
23:DB:20:C:H2'	23:DB:21:A:H8	1.64	0.63
23:DB:850:U:H2'	23:DB:851:C:C6	2.34	0.63
23:DB:946:C:H2'	23:DB:947:A:C8	2.33	0.63
24:DI:27:LEU:HD12	24:DI:32:VAL:HG11	1.79	0.63
24:DI:41:PHE:O	24:DI:45:THR:HG23	1.98	0.63
41:DJ:98:GLU:HB3	41:DJ:124:VAL:HG21	1.79	0.63
23:DB:811:U:H2'	37:DL:21:ARG:HA	1.79	0.63
49:DR:24:LYS:HA	49:DR:94:THR:OG1	1.97	0.63
45:DS:60:HIS:O	45:DS:61:ASN:HB2	1.98	0.63
52:DW:49:ASN:HB3	52:DW:60:ALA:HA	1.80	0.63
1:AA:1047:G:H21	1:AA:1215:G:H5''	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1409:C:H4'	23:BB:1914:C:N4	2.12	0.63
1:AA:26:A:H61	1:AA:558:G:H1'	1.64	0.63
1:AA:947:G:H2'	1:AA:948:C:C6	2.33	0.63
7:AH:58:LEU:HD22	7:AH:60:LEU:HB2	1.80	0.63
8:AI:17:ARG:HE	8:AI:65:THR:HG21	1.62	0.63
8:AI:98:ARG:HA	8:AI:103:VAL:HG22	1.80	0.63
13:AP:20:VAL:HG23	13:AP:35:ARG:HA	1.78	0.63
19:AU:13:VAL:HG22	19:AU:15:LEU:HD23	1.81	0.63
19:AU:36:PHE:HB3	19:AU:40:PRO:CD	2.26	0.63
23:BB:1028:A:H2'	23:BB:1029:A:C8	2.34	0.63
23:BB:129:C:H2'	23:BB:130:C:C6	2.34	0.63
23:BB:3:U:O2'	23:BB:4:U:H6	1.81	0.63
25:BC:116:GLN:HG2	25:BC:117:SER:H	1.63	0.63
25:BC:93:VAL:O	25:BC:94:LEU:HB3	1.98	0.63
48:BG:84:LYS:HG3	48:BG:131:VAL:C	2.19	0.63
41:BJ:13:ARG:O	41:BJ:14:ASP:HB2	1.98	0.63
41:BJ:29:ALA:HA	41:BJ:32:LEU:HD12	1.80	0.63
44:BQ:57:ARG:NH1	44:BQ:61:ILE:HD11	2.13	0.63
50:BT:29:THR:CA	50:BT:86:THR:HA	2.26	0.63
35:BV:68:LYS:HD3	35:BV:68:LYS:H	1.62	0.63
2:CC:26:LYS:HG3	2:CC:27:GLU:N	2.10	0.63
4:CE:92:ARG:HB3	4:CE:92:ARG:HH11	1.63	0.63
5:CF:3:HIS:CE1	5:CF:95:ALA:H	2.17	0.63
8:CI:41:GLU:O	8:CI:44:ARG:HG2	1.98	0.63
9:CJ:53:ILE:HG13	21:CN:84:ARG:NE	2.12	0.63
23:DB:2734:A:H2'	23:DB:2735:G:H5'	1.81	0.63
23:DB:857:G:O2'	23:DB:858:G:H5'	1.98	0.63
26:DD:178:VAL:HB	26:DD:188:LEU:HB2	1.81	0.63
42:DN:33:ILE:HD11	42:DN:112:TYR:HD1	1.62	0.63
28:DP:59:THR:H	28:DP:72:VAL:HA	1.63	0.63
45:DS:46:LEU:O	45:DS:50:VAL:HG23	1.98	0.63
1:AA:1095:U:H2'	1:AA:1096:C:C6	2.34	0.63
1:AA:1458:G:H2'	1:AA:1459:G:C8	2.33	0.63
1:AA:22:G:H2'	1:AA:23:C:C6	2.33	0.63
1:AA:423:G:H2'	1:AA:424:G:O4'	1.97	0.63
1:AA:430:A:OP1	3:AD:8:LEU:HB2	1.99	0.63
1:AA:674:G:H2'	1:AA:675:A:C8	2.32	0.63
13:AP:66:THR:HG22	13:AP:67:ILE:N	2.10	0.63
23:BB:1406:U:H2'	23:BB:1407:G:H8	1.63	0.63
23:BB:609:A:H2'	23:BB:610:C:O4'	1.99	0.63
26:BD:33:ARG:HH11	26:BD:51:THR:CG2	2.11	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:8:LYS:HB2	26:BD:201:LEU:HD11	1.79	0.63
29:BE:146:VAL:HA	29:BE:185:LYS:O	1.97	0.63
43:BO:90:VAL:HG22	43:BO:115:LEU:HD12	1.80	0.63
45:BS:15:GLN:HA	45:BS:18:ARG:HG2	1.81	0.63
1:CA:26:A:N6	1:CA:558:G:H1'	2.13	0.63
2:CC:51:VAL:HA	2:CC:69:THR:HA	1.81	0.63
1:CA:1329:A:O2'	12:CM:23:GLY:HA2	1.98	0.63
16:CS:52:ASN:HD22	16:CS:76:THR:HA	1.62	0.63
23:DB:2243:U:O2'	23:DB:2244:U:H5'	1.99	0.63
23:DB:3:U:H2'	23:DB:4:U:C6	2.33	0.63
23:DB:2579:C:O2'	26:DD:136:ASN:HA	1.99	0.63
1:AA:1324:A:H2'	1:AA:1325:C:C6	2.34	0.63
1:AA:209:U:H5'	1:AA:210:C:H5	1.63	0.63
1:AA:26:A:N6	1:AA:558:G:H1'	2.13	0.63
1:AA:270:A:H2'	1:AA:271:C:C6	2.33	0.63
1:AA:441:A:H61	1:AA:493:A:N6	1.97	0.63
1:AA:86:G:N2	1:AA:87:C:N4	2.47	0.63
2:AC:153:SER:O	2:AC:164:THR:HA	1.98	0.63
11:AL:98:ARG:HA	11:AL:103:CYS:SG	2.39	0.63
12:AM:9:PRO:HG2	12:AM:44:ILE:HG21	1.81	0.63
21:AN:61:ASN:HB3	21:AN:72:PHE:CE2	2.33	0.63
22:BA:32:U:H1'	22:BA:52:A:N7	2.13	0.63
23:BB:1812:U:H2'	23:BB:1813:G:C8	2.32	0.63
23:BB:2893:A:H5''	23:BB:2894:G:H5'	1.79	0.63
23:BB:2313:C:H4'	47:BF:87:LYS:HB3	1.81	0.63
38:BM:68:PHE:CD1	38:BM:69:PRO:HD2	2.33	0.63
44:BQ:87:VAL:HG21	49:BR:52:PRO:HG3	1.79	0.63
52:BW:49:ASN:OD1	52:BW:81:ILE:HG12	1.98	0.63
1:CA:1014:A:H2'	1:CA:1015:G:O4'	1.99	0.63
1:CA:845:A:H5''	1:CA:846:G:C8	2.33	0.63
4:CE:131:ASN:HD22	4:CE:134:ASN:H	1.47	0.63
6:CG:68:VAL:O	6:CG:137:ARG:HG3	1.99	0.63
10:CK:62:ALA:HB1	10:CK:95:THR:HB	1.80	0.63
12:CM:65:GLU:H	12:CM:68:LEU:HD12	1.64	0.63
1:CA:1320:C:OP2	16:CS:2:ARG:HG3	1.98	0.63
23:DB:1173:U:H2'	23:DB:1174:U:N1	2.13	0.63
23:DB:1248:G:OP1	44:DQ:1:ALA:HB3	1.99	0.63
23:DB:137:U:H2'	23:DB:138:U:O4'	1.99	0.63
23:DB:1487:U:H2'	23:DB:1488:C:H6	1.64	0.63
23:DB:1548:A:H2'	23:DB:1549:A:H8	1.63	0.63
23:DB:2461:A:H2'	23:DB:2462:C:C6	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:37:C:O2'	29:DE:45:ALA:HA	1.99	0.63
23:DB:5:A:H2'	23:DB:6:A:C8	2.33	0.63
23:DB:62:U:H2'	23:DB:63:A:O4'	1.99	0.63
25:DC:116:GLN:HG2	25:DC:117:SER:H	1.63	0.63
26:DD:33:ARG:HD3	26:DD:74:GLU:OE1	1.99	0.63
47:DF:34:THR:HA	47:DF:89:THR:HG22	1.79	0.63
37:DL:143:GLU:HG2	37:DL:144:GLU:N	2.12	0.63
37:DL:124:GLY:H	37:DL:143:GLU:HG3	1.63	0.63
28:DP:20:ARG:O	28:DP:46:VAL:HG21	1.99	0.63
50:DT:7:LEU:C	50:DT:9:LYS:HD2	2.18	0.63
52:DW:49:ASN:O	52:DW:50:VAL:HG13	1.99	0.63
1:AA:1137:C:O2'	1:AA:1138:G:H5''	1.99	0.63
1:AA:1162:C:H2'	1:AA:1163:A:H8	1.63	0.63
2:AC:35:ASP:O	2:AC:39:ARG:HG3	1.99	0.63
4:AE:132:PRO:O	4:AE:136:VAL:HG12	1.98	0.63
6:AG:50:ALA:CB	6:AG:57:GLU:HG3	2.29	0.63
7:AH:73:SER:O	7:AH:128:VAL:HA	1.99	0.63
8:AI:75:ALA:HA	8:AI:78:ILE:HD12	1.80	0.63
9:AJ:51:VAL:HG23	21:AN:80:ARG:HD2	1.80	0.63
9:AJ:8:ILE:HG22	9:AJ:8:ILE:O	1.99	0.63
23:BB:1189:A:H2'	23:BB:1190:G:O4'	1.99	0.63
23:BB:150:U:H2'	23:BB:151:C:H6	1.63	0.63
23:BB:1913:A:H1'	23:BB:1914:C:OP1	1.99	0.63
23:BB:2734:A:H2'	23:BB:2735:G:H5'	1.80	0.63
25:BC:244:VAL:HB	25:BC:249:VAL:N	2.08	0.63
26:BD:4:LEU:HG	26:BD:29:VAL:HG11	1.80	0.63
47:BF:36:ASN:HA	47:BF:87:LYS:HA	1.80	0.63
45:BS:60:HIS:O	45:BS:61:ASN:HB2	1.97	0.63
1:CA:1078:U:H4'	4:CE:137:ARG:NH1	2.13	0.63
1:CA:1234:C:O2'	1:CA:1235:U:H5'	1.99	0.63
1:CA:1278:G:H4'	1:CA:1279:G:O5'	1.99	0.63
1:CA:1356:G:O2'	1:CA:1357:A:H5'	1.98	0.63
1:CA:677:U:H2'	1:CA:678:U:H6	1.62	0.63
5:CF:29:ILE:HG21	5:CF:64:VAL:HG11	1.81	0.63
10:CK:52:ARG:HH11	10:CK:52:ARG:HB3	1.62	0.63
22:DA:43:C:H2'	22:DA:44:G:H5''	1.80	0.63
23:DB:1060:U:O4	23:DB:1088:A:N6	2.31	0.63
23:DB:1198:U:H2'	23:DB:1199:U:C6	2.34	0.63
1:CA:1409:C:H1'	23:DB:1913:A:N6	2.14	0.63
23:DB:2516:A:O2'	23:DB:2517:C:H5'	1.98	0.63
23:DB:2592:G:H2'	23:DB:2593:U:O4'	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:6:A:H2'	23:DB:7:G:H8	1.64	0.63
24:DI:121:ILE:HD13	24:DI:121:ILE:N	2.13	0.63
44:DQ:4:LYS:NZ	44:DQ:7:VAL:HG22	2.14	0.63
46:DU:27:VAL:HG23	46:DU:33:VAL:HG12	1.79	0.63
39:DX:23:ARG:O	39:DX:27:ASN:HB2	1.99	0.63
1:AA:1081:A:H2'	1:AA:1082:A:C8	2.34	0.63
1:AA:411:A:H2'	1:AA:412:A:N3	2.14	0.63
1:AA:414:A:H2'	1:AA:415:A:O4'	1.98	0.63
8:AI:11:ARG:O	8:AI:14:SER:HB2	1.99	0.63
9:AJ:6:ILE:HB	9:AJ:76:ILE:HG13	1.79	0.63
23:BB:2008:C:H2'	23:BB:2009:A:C8	2.33	0.63
23:BB:2341:G:H2'	23:BB:2342:C:C6	2.33	0.63
23:BB:2547:A:H2'	23:BB:2548:U:C6	2.34	0.63
38:BM:105:MET:HB2	38:BM:117:PHE:HZ	1.63	0.63
28:BP:88:ARG:HB2	28:BP:112:ARG:NH1	2.13	0.63
23:BB:1599:U:OP1	50:BT:40:LYS:HB3	1.99	0.63
23:BB:2269:G:H4'	52:BW:19:ARG:HH11	1.64	0.63
1:CA:312:C:H2'	1:CA:313:A:C8	2.32	0.63
1:CA:414:A:H2'	1:CA:415:A:O4'	1.99	0.63
1:CA:98:A:H2'	1:CA:99:C:C6	2.34	0.63
18:CB:127:LYS:HD2	18:CB:128:LEU:N	2.13	0.63
6:CG:19:SER:CB	6:CG:22:LEU:HB2	2.28	0.63
6:CG:2:ARG:HB3	6:CG:2:ARG:NH1	2.14	0.63
11:CL:34:THR:HB	11:CL:53:ARG:HB3	1.80	0.63
15:CR:60:ARG:HA	15:CR:63:TYR:HD2	1.62	0.63
23:DB:1725:U:H2'	23:DB:1726:C:C6	2.34	0.63
23:DB:2031:A:H8	23:DB:2031:A:H5'	1.62	0.63
23:DB:2195:U:O2'	23:DB:2196:C:H5'	1.99	0.63
23:DB:705:A:N6	23:DB:726:G:H1'	2.14	0.63
40:DH:113:SER:O	40:DH:115:VAL:N	2.32	0.63
38:DM:66:ARG:NE	38:DM:101:VAL:HG11	2.14	0.63
1:AA:1165:U:H2'	1:AA:1166:G:O4'	1.98	0.63
3:AD:84:ASN:ND2	3:AD:86:GLY:H	1.97	0.63
6:AG:149:ALA:HB1	10:AK:58:THR:HB	1.79	0.63
20:AO:57:LEU:O	20:AO:60:VAL:HB	1.98	0.63
34:B3:61:LEU:N	34:B3:62:PRO:HD3	2.14	0.63
23:BB:1060:U:O4	23:BB:1088:A:N6	2.32	0.63
23:BB:1251:C:H2'	44:BQ:5:ARG:NH1	2.11	0.63
23:BB:2021:C:OP1	31:B0:8:THR:HG21	1.98	0.63
23:BB:20:C:H2'	23:BB:21:A:H8	1.62	0.63
23:BB:560:C:H3'	23:BB:561:G:C8	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:68:PHE:HB3	26:BD:73:VAL:HG23	1.81	0.63
24:BI:105:LEU:HD11	24:BI:139:VAL:HG11	1.80	0.63
45:BS:84:ARG:HB3	45:BS:96:ILE:HG23	1.81	0.63
1:CA:979:C:H1'	1:CA:1317:C:N4	2.14	0.63
2:CC:72:PRO:O	2:CC:76:ILE:HG12	1.99	0.63
9:CJ:9:ARG:HB2	9:CJ:99:GLN:HB3	1.81	0.63
32:D4:10:LEU:HD12	32:D4:33:HIS:HA	1.80	0.63
32:D4:5:ALA:CA	32:D4:37:GLN:HE22	2.12	0.63
23:DB:324:A:H2'	23:DB:325:G:O4'	1.98	0.63
25:DC:202:ARG:O	25:DC:203:VAL:HB	1.99	0.63
29:DE:165:HIS:HD2	29:DE:166:LYS:HD3	1.62	0.63
23:DB:674:G:H5''	29:DE:71:GLY:H	1.63	0.63
40:DH:41:LYS:O	40:DH:44:ILE:HG12	1.98	0.63
24:DI:72:THR:HG22	24:DI:115:ASP:OD2	1.98	0.63
38:DM:78:LEU:H	38:DM:78:LEU:HD12	1.64	0.63
52:DW:37:VAL:HG12	52:DW:38:ARG:N	2.10	0.63
1:AA:1164:G:H2'	1:AA:1165:U:H6	1.64	0.62
1:AA:32:A:H2'	1:AA:33:A:C8	2.34	0.62
1:AA:676:A:O2'	1:AA:677:U:H5'	1.99	0.62
3:AD:29:THR:HB	3:AD:30:LYS:NZ	2.14	0.62
16:AS:52:ASN:ND2	16:AS:53:GLY:H	1.97	0.62
33:B1:8:ILE:CD1	33:B1:51:ALA:HA	2.29	0.62
23:BB:1081:U:H4'	24:BI:123:ALA:HB1	1.80	0.62
23:BB:143:C:O4'	50:BT:2:ILE:HB	1.99	0.62
23:BB:191:A:H2'	23:BB:192:C:H6	1.64	0.62
23:BB:2425:A:H5'	23:BB:2427:C:O4'	1.99	0.62
23:BB:275:C:C2'	23:BB:276:U:H5'	2.29	0.62
23:BB:876:C:H41	23:BB:901:C:H41	1.47	0.62
23:BB:974:G:OP2	49:BR:78:ARG:HD3	1.99	0.62
25:BC:130:PRO:HA	25:BC:188:ARG:HA	1.81	0.62
40:BH:82:SER:H	40:BH:146:VAL:HG13	1.63	0.62
37:BL:111:ILE:HG22	37:BL:112:LEU:H	1.64	0.62
22:BA:98:G:H1	35:BV:14:LYS:HB2	1.63	0.62
1:CA:1392:G:O2'	1:CA:1393:U:H5'	1.99	0.62
1:CA:865:A:H2'	1:CA:866:C:C6	2.34	0.62
3:CD:25:ARG:NH1	3:CD:26:ALA:HB2	2.14	0.62
8:CI:43:ALA:O	8:CI:46:VAL:HG22	1.99	0.62
8:CI:22:PRO:HA	8:CI:60:LEU:HB3	1.81	0.62
21:CN:98:ALA:HB1	21:CN:100:TRP:HZ3	1.64	0.62
23:DB:150:U:H2'	23:DB:151:C:C6	2.34	0.62
23:DB:2537:U:H2'	23:DB:2538:C:C6	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:DG:15:ASP:HB2	48:DG:26:LYS:CD	2.29	0.62
41:DJ:136:GLN:N	41:DJ:137:PRO:HD3	2.14	0.62
50:DT:29:THR:CA	50:DT:86:THR:HA	2.27	0.62
35:DV:80:HIS:CD2	35:DV:83:LYS:HB2	2.34	0.62
8:AI:43:ALA:HA	8:AI:46:VAL:HG13	1.79	0.62
21:AN:43:ALA:HA	21:AN:46:LYS:HG2	1.81	0.62
17:AT:43:LYS:H	17:AT:43:LYS:HD3	1.62	0.62
32:B4:5:ALA:CA	32:B4:37:GLN:HE22	2.12	0.62
23:BB:1725:U:H2'	23:BB:1726:C:C6	2.34	0.62
23:BB:2031:A:C6	23:BB:2498:C:H1'	2.34	0.62
23:BB:414:C:H2'	23:BB:415:A:C8	2.34	0.62
26:BD:168:GLU:HG3	26:BD:170:VAL:HG13	1.80	0.62
40:BH:81:ALA:HA	40:BH:146:VAL:HA	1.81	0.62
27:BK:110:GLU:HA	27:BK:113:MET:HG3	1.81	0.62
35:BV:75:GLN:HG2	35:BV:92:VAL:HG23	1.81	0.62
1:CA:1320:C:N3	16:CS:35:ARG:HD3	2.14	0.62
1:CA:967:C:H1'	8:CI:129:ARG:NH1	2.14	0.62
21:CN:68:ARG:HH11	21:CN:70:HIS:HB2	1.63	0.62
23:DB:1640:A:H2'	23:DB:1641:A:C8	2.34	0.62
23:DB:1843:C:H5''	25:DC:250:GLN:NE2	2.15	0.62
23:DB:2241:A:H2'	23:DB:2242:G:C8	2.33	0.62
23:DB:2671:G:H2'	23:DB:2672:U:C6	2.35	0.62
23:DB:282:A:H2'	23:DB:283:G:O4'	1.98	0.62
23:DB:2898:U:H2'	23:DB:2899:A:C8	2.34	0.62
26:DD:15:PHE:N	28:DP:11:GLN:HE22	1.97	0.62
29:DE:52:VAL:HG12	29:DE:53:THR:N	2.13	0.62
48:DG:86:LEU:HD23	48:DG:162:ARG:O	1.99	0.62
48:DG:24:THR:HA	48:DG:33:THR:O	1.99	0.62
44:DQ:89:ILE:HG21	49:DR:11:GLN:NE2	2.12	0.62
45:DS:10:ALA:HB3	45:DS:101:SER:OG	1.97	0.62
46:DU:26:ASN:ND2	46:DU:34:ILE:HD12	2.14	0.62
52:DW:23:LYS:HD2	52:DW:24:ARG:HB3	1.81	0.62
1:AA:1240:U:H4'	1:AA:1241:G:OP2	1.98	0.62
1:AA:1492:A:N7	1:AA:1493:A:H1'	2.14	0.62
1:AA:673:A:H2'	1:AA:674:G:C8	2.34	0.62
1:AA:817:C:H1'	1:AA:819:A:H5'	1.81	0.62
8:AI:26:LYS:N	8:AI:61:ASP:HB3	2.13	0.62
23:BB:1470:A:H3'	23:BB:1471:G:H8	1.62	0.62
23:BB:1845:G:O2'	23:BB:1846:G:H5'	1.99	0.62
23:BB:2575:C:H5''	26:BD:149:ASN:HD22	1.63	0.62
23:BB:2678:C:H2'	23:BB:2679:A:H8	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:165:HIS:HD2	29:BE:166:LYS:HD3	1.63	0.62
38:BM:19:GLY:N	38:BM:38:ARG:HH12	1.98	0.62
52:BW:37:VAL:HG13	52:BW:55:ASP:O	1.99	0.62
52:BW:76:ARG:HA	52:BW:76:ARG:NH2	2.14	0.62
30:BY:7:THR:HG23	30:BY:34:THR:OG1	1.99	0.62
1:CA:865:A:H5'	1:CA:1078:U:O4	1.99	0.62
1:CA:1365:G:H2'	1:CA:1366:C:H6	1.64	0.62
6:CG:75:LYS:HZ2	6:CG:76:SER:N	1.92	0.62
7:CH:44:PHE:HE2	7:CH:100:ILE:HG12	1.64	0.62
1:CA:1060:U:H4'	9:CJ:54:SER:CB	2.27	0.62
14:CQ:18:LYS:H	14:CQ:50:ASN:HD21	1.46	0.62
17:CT:43:LYS:HB3	17:CT:86:ALA:HB1	1.80	0.62
17:CT:74:HIS:O	17:CT:78:LEU:HG	1.99	0.62
23:DB:1196:C:H2'	23:DB:1197:G:C8	2.34	0.62
23:DB:1774:C:O2	23:DB:1774:C:H2'	1.97	0.62
23:DB:2031:A:C6	23:DB:2498:C:H1'	2.35	0.62
26:DD:62:LYS:HB2	26:DD:63:PRO:HD3	1.79	0.62
23:DB:1652:A:N6	42:DN:11:ASN:HD21	1.96	0.62
43:DO:90:VAL:HG22	43:DO:115:LEU:HD12	1.80	0.62
35:DV:53:LYS:HZ2	35:DV:54:ALA:H	1.46	0.62
51:DZ:35:SER:HA	51:DZ:50:ARG:HA	1.80	0.62
5:AF:3:HIS:CE1	5:AF:95:ALA:H	2.17	0.62
21:AN:76:PHE:O	21:AN:78:LEU:HD13	1.99	0.62
32:B4:22:VAL:HG11	32:B4:36:ARG:HD2	1.81	0.62
23:BB:2615:U:C2	31:B0:3:GLN:HA	2.35	0.62
25:BC:141:HIS:HB3	25:BC:190:THR:OG1	2.00	0.62
47:BF:84:ILE:HG23	47:BF:85:GLY:N	2.12	0.62
52:BW:17:ALA:CA	52:BW:35:ILE:HG23	2.22	0.62
51:BZ:45:ARG:NE	51:BZ:47:VAL:HG12	2.03	0.62
1:CA:1492:A:H2'	1:CA:1492:A:N3	2.13	0.62
1:CA:925:G:O2'	1:CA:926:G:H5''	1.98	0.62
2:CC:133:MET:HE3	2:CC:152:VAL:HG23	1.82	0.62
3:CD:121:ALA:O	3:CD:122:ILE:HD13	1.98	0.62
12:CM:33:LEU:HD13	12:CM:39:ALA:O	1.98	0.62
21:CN:50:LEU:HD23	21:CN:51:PRO:HD3	1.81	0.62
19:CU:40:PRO:HA	19:CU:44:ARG:HD2	1.80	0.62
23:DB:347:A:H2'	23:DB:348:A:C8	2.35	0.62
48:DG:95:ALA:HB2	48:DG:130:ILE:HD11	1.81	0.62
23:DB:1080:A:O2'	24:DI:126:ARG:HD2	1.99	0.62
46:DU:26:ASN:N	46:DU:26:ASN:HD22	1.96	0.62
1:AA:1182:G:H4'	1:AA:1183:U:H5'	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1218:C:H2'	1:AA:1219:A:C8	2.34	0.62
1:AA:1271:A:H5'	1:AA:1314:C:H5''	1.80	0.62
1:AA:918:A:H2'	1:AA:919:A:C8	2.34	0.62
2:AC:152:VAL:O	2:AC:164:THR:HG23	1.99	0.62
6:AG:52:ARG:HH22	6:AG:120:ALA:HB1	1.65	0.62
8:AI:80:HIS:NE2	8:AI:84:ARG:HD3	2.14	0.62
16:AS:43:MET:O	16:AS:46:LEU:HB2	2.00	0.62
10:AK:92:ARG:NH2	19:AU:19:LYS:HG2	2.15	0.62
23:BB:2348:U:OP1	34:B3:37:THR:HG21	2.00	0.62
32:B4:16:ILE:HA	32:B4:25:VAL:HA	1.80	0.62
32:B4:10:LEU:HD12	32:B4:33:HIS:HA	1.80	0.62
23:BB:1993:U:H4'	26:BD:133:THR:CG2	2.29	0.62
23:BB:2810:A:H2'	23:BB:2811:G:O4'	1.99	0.62
23:BB:364:C:H2'	23:BB:365:U:H6	1.61	0.62
23:BB:6:A:H2'	23:BB:7:G:C8	2.35	0.62
26:BD:149:ASN:O	26:BD:152:PRO:HD2	1.99	0.62
26:BD:178:VAL:HB	26:BD:188:LEU:HB2	1.81	0.62
47:BF:109:ARG:HB2	47:BF:135:ILE:HD12	1.82	0.62
48:BG:144:ALA:HB1	48:BG:163:TYR:HE1	1.65	0.62
39:BX:13:GLU:O	39:BX:17:GLU:HG3	1.99	0.62
30:BY:9:THR:HB	30:BY:53:MET:O	1.99	0.62
6:CG:26:VAL:HB	6:CG:42:VAL:HG21	1.81	0.62
7:CH:23:ALA:HA	7:CH:62:LEU:H	1.64	0.62
16:CS:44:ILE:HA	16:CS:61:VAL:HB	1.81	0.62
22:DA:87:U:H2'	22:DA:88:C:O5'	1.98	0.62
23:DB:135:U:H2'	23:DB:136:G:C8	2.34	0.62
23:DB:2866:U:H4'	23:DB:2867:G:H4'	1.81	0.62
23:DB:2880:C:O4'	42:DN:91:ALA:HB3	2.00	0.62
25:DC:141:HIS:HB3	25:DC:190:THR:OG1	2.00	0.62
47:DF:101:ARG:NH1	47:DF:138:PRO:HB2	2.15	0.62
41:DJ:13:ARG:O	41:DJ:14:ASP:HB2	1.99	0.62
37:DL:80:SER:H	37:DL:113:ALA:HB3	1.63	0.62
45:DS:17:VAL:C	45:DS:19:LEU:H	2.00	0.62
1:AA:74:A:H3'	1:AA:75:G:H5''	1.81	0.62
1:AA:922:G:H2'	1:AA:923:A:H8	1.63	0.62
18:AB:112:ARG:HH21	18:AB:116:LEU:HD11	1.64	0.62
9:AJ:7:ARG:HA	9:AJ:75:ASP:OD2	2.00	0.62
12:AM:89:ARG:NH1	12:AM:94:LEU:HB3	2.15	0.62
12:AM:9:PRO:O	12:AM:44:ILE:HB	1.99	0.62
15:AR:31:TYR:O	15:AR:39:VAL:HG22	1.98	0.62
23:BB:946:C:H2'	23:BB:947:A:C8	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:136:GLN:O	29:BE:139:LYS:HG2	2.00	0.62
41:BJ:136:GLN:N	41:BJ:137:PRO:HD3	2.15	0.62
28:BP:63:ILE:HA	28:BP:68:GLY:HA2	1.80	0.62
52:BW:37:VAL:HG12	52:BW:38:ARG:N	2.12	0.62
1:CA:451:A:H5'	13:CP:70:ARG:NH2	2.14	0.62
1:CA:812:G:H2'	1:CA:812:G:N3	2.14	0.62
11:CL:31:GLY:HA3	11:CL:54:VAL:CG1	2.30	0.62
13:CP:67:ILE:HG12	13:CP:72:ALA:HB2	1.82	0.62
22:DA:49:C:H2'	22:DA:50:A:C8	2.33	0.62
23:DB:1577:C:H2'	23:DB:1578:U:C6	2.34	0.62
23:DB:1594:U:H2'	23:DB:1595:C:C6	2.34	0.62
23:DB:1796:U:H2'	23:DB:1797:G:C8	2.33	0.62
23:DB:2586:U:H2'	23:DB:2587:A:C8	2.34	0.62
23:DB:2626:C:H2'	23:DB:2627:G:C8	2.35	0.62
23:DB:608:A:H2'	23:DB:609:A:C8	2.34	0.62
30:DY:26:LEU:HD12	30:DY:28:LEU:HD11	1.81	0.62
1:AA:1492:A:H2'	23:BB:1913:A:N6	2.14	0.62
1:AA:194:C:O2'	1:AA:195:A:H5'	1.99	0.62
1:AA:57:G:H2'	1:AA:58:C:C6	2.34	0.62
1:AA:83:C:H4'	1:AA:84:U:C5	2.34	0.62
4:AE:92:ARG:HH11	4:AE:92:ARG:HB3	1.63	0.62
8:AI:98:ARG:NE	8:AI:103:VAL:HG21	2.15	0.62
22:BA:43:C:H2'	22:BA:44:G:H5''	1.80	0.62
23:BB:1458:U:H5''	23:BB:1459:G:OP1	2.00	0.62
23:BB:2284:A:OP2	33:B1:5:ARG:HG3	2.00	0.62
23:BB:2286:G:H4'	23:BB:2287:A:O4'	2.00	0.62
23:BB:2743:U:H2'	23:BB:2744:G:O4'	2.00	0.62
26:BD:33:ARG:HH11	26:BD:51:THR:HG22	1.64	0.62
26:BD:56:LYS:HD2	26:BD:59:ARG:HG3	1.80	0.62
47:BF:77:LYS:HD2	47:BF:79:ARG:HD3	1.81	0.62
48:BG:145:ALA:O	48:BG:148:ARG:HG3	2.00	0.62
40:BH:116:ARG:H	40:BH:130:VAL:HG12	1.64	0.62
28:BP:20:ARG:O	28:BP:46:VAL:HG21	1.99	0.62
44:BQ:91:ARG:HB2	44:BQ:94:LEU:HD23	1.82	0.62
1:CA:1278:G:H4'	1:CA:1279:G:H5'	1.81	0.62
1:CA:270:A:H2'	1:CA:271:C:C6	2.34	0.62
6:CG:14:ASP:N	6:CG:19:SER:H	1.93	0.62
1:CA:716:A:N3	10:CK:119:GLY:HA2	2.15	0.62
22:DA:91:C:H2'	22:DA:92:C:C6	2.34	0.62
23:DB:1810:A:H2'	23:DB:1811:G:O4'	1.99	0.62
23:DB:2892:G:H5''	23:DB:2894:G:N2	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DM:68:PHE:CD1	38:DM:69:PRO:HD2	2.34	0.62
45:DS:15:GLN:HA	45:DS:18:ARG:HG2	1.81	0.62
1:AA:1114:C:H2'	1:AA:1115:U:O4'	1.99	0.62
1:AA:900:A:H2'	1:AA:901:A:C8	2.34	0.62
3:AD:172:VAL:HA	3:AD:179:GLY:HA2	1.80	0.62
8:AI:74:GLN:O	8:AI:78:ILE:HG13	1.99	0.62
34:B3:39:ARG:O	34:B3:43:LEU:HG	2.00	0.62
23:BB:1551:A:C3'	23:BB:1552:A:H5''	2.29	0.62
23:BB:2592:G:H2'	23:BB:2593:U:O4'	2.00	0.62
23:BB:2677:G:H2'	23:BB:2678:C:C6	2.34	0.62
23:BB:303:G:H2'	23:BB:304:U:C6	2.35	0.62
23:BB:784:G:O2'	23:BB:785:G:H5''	2.00	0.62
23:BB:83:A:N6	23:BB:101:A:H5'	2.15	0.62
29:BE:132:LYS:O	29:BE:135:ALA:HB3	1.99	0.62
47:BF:78:ILE:HA	47:BF:82:TYR:CG	2.34	0.62
48:BG:86:LEU:HD23	48:BG:162:ARG:O	1.99	0.62
24:BI:85:ILE:HD13	24:BI:137:LEU:HD21	1.81	0.62
1:CA:1262:C:N4	1:CA:1273:C:H42	1.98	0.62
1:CA:517:G:H22	1:CA:533:A:P	2.23	0.62
1:CA:729:A:H2'	1:CA:730:G:H8	1.65	0.62
6:CG:23:ALA:O	6:CG:26:VAL:HG22	1.99	0.62
21:CN:50:LEU:H	21:CN:51:PRO:CD	2.11	0.62
21:CN:79:SER:OG	21:CN:82:LYS:HG2	2.00	0.62
13:CP:66:THR:HG22	13:CP:67:ILE:N	2.13	0.62
23:DB:140:C:H5'	23:DB:141:G:C6	2.34	0.62
23:DB:609:A:H2'	23:DB:610:C:O4'	1.99	0.62
23:DB:6:A:H2'	23:DB:7:G:C8	2.35	0.62
29:DE:136:GLN:O	29:DE:139:LYS:HG2	1.99	0.62
50:DT:2:ILE:HG12	50:DT:3:ARG:HD3	1.82	0.62
52:DW:49:ASN:OD1	52:DW:81:ILE:HG12	1.98	0.62
1:AA:1053:G:O6	1:AA:1200:C:H5''	1.99	0.62
1:AA:1448:C:H2'	1:AA:1449:C:H6	1.64	0.62
1:AA:996:A:H2'	1:AA:997:U:C6	2.35	0.62
12:AM:84:CYS:O	12:AM:88:LEU:HG	1.98	0.62
31:B0:9:ARG:O	31:B0:12:ARG:HB3	1.99	0.62
32:B4:24:ARG:HG2	32:B4:36:ARG:HG3	1.82	0.62
23:BB:116:C:O2'	23:BB:117:G:H5'	2.00	0.62
23:BB:1196:C:H2'	23:BB:1197:G:C8	2.35	0.62
23:BB:2064:C:H2'	23:BB:2065:C:H6	1.64	0.62
23:BB:2455:G:H2'	23:BB:2456:C:C6	2.35	0.62
23:BB:320:A:H4'	23:BB:322:A:N7	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:16:VAL:HB	25:BC:203:VAL:HG11	1.82	0.62
29:BE:178:VAL:O	29:BE:181:ILE:HG22	2.00	0.62
50:BT:2:ILE:HD11	50:BT:3:ARG:CZ	2.30	0.62
1:CA:474:G:H2'	1:CA:475:C:C6	2.34	0.62
1:CA:624:C:H2'	1:CA:625:U:H6	1.64	0.62
1:CA:918:A:H2'	1:CA:919:A:C8	2.35	0.62
20:CO:11:ILE:HA	20:CO:14:GLU:OE1	1.99	0.62
23:DB:1115:G:H2'	23:DB:1116:G:C8	2.35	0.62
47:DF:177:ARG:HA	47:DF:177:ARG:NH1	2.14	0.62
37:DL:95:LEU:HB2	37:DL:101:ILE:HG13	1.81	0.62
28:DP:21:PRO:O	28:DP:46:VAL:HG23	2.00	0.62
1:AA:382:A:H2'	1:AA:383:A:C8	2.34	0.62
1:AA:451:A:H5'	13:AP:70:ARG:NH2	2.15	0.62
18:AB:169:HIS:HA	18:AB:172:ILE:HD12	1.82	0.62
18:AB:27:LYS:HA	18:AB:30:ILE:HD12	1.80	0.62
18:AB:45:THR:HG22	18:AB:49:PHE:CE1	2.35	0.62
6:AG:4:ARG:HE	6:AG:6:ILE:HG23	1.65	0.62
11:AL:41:PRO:HB3	11:AL:49:ARG:NH1	2.15	0.62
9:AJ:65:TYR:C	21:AN:98:ALA:HB2	2.21	0.62
34:B3:54:LEU:O	34:B3:57:VAL:HG23	2.00	0.62
23:BB:1220:G:H2'	23:BB:1221:C:C6	2.35	0.62
23:BB:1354:A:H2'	23:BB:1355:G:O4'	2.00	0.62
23:BB:1406:U:H2'	23:BB:1407:G:C8	2.35	0.62
23:BB:1487:U:H2'	23:BB:1488:C:H6	1.64	0.62
23:BB:2866:U:H4'	23:BB:2867:G:H4'	1.81	0.62
23:BB:533:G:H2'	23:BB:534:U:C6	2.34	0.62
23:BB:608:A:H2'	23:BB:609:A:C8	2.35	0.62
25:BC:43:ASN:HB2	25:BC:49:THR:HG23	1.82	0.62
25:BC:92:LEU:HD12	25:BC:93:VAL:H	1.64	0.62
27:BK:87:LEU:HB2	27:BK:93:GLN:O	2.00	0.62
46:BU:42:LYS:HG3	46:BU:57:ILE:HG21	1.82	0.62
1:CA:1097:C:H2'	1:CA:1098:C:C6	2.35	0.62
1:CA:57:G:H2'	1:CA:58:C:C6	2.35	0.62
1:CA:764:C:C2'	1:CA:765:G:H5'	2.30	0.62
18:CB:40:ILE:HD13	18:CB:201:GLY:CA	2.26	0.62
2:CC:75:VAL:O	2:CC:82:ASP:HB2	2.00	0.62
1:CA:1374:A:H4'	6:CG:27:ASN:OD1	2.00	0.62
8:CI:20:ILE:HG13	8:CI:62:LEU:HD12	1.81	0.62
32:D4:16:ILE:HA	32:D4:25:VAL:HA	1.81	0.62
23:DB:150:U:H2'	23:DB:151:C:H6	1.64	0.62
23:DB:1821:A:H2'	23:DB:1822:C:H6	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:532:A:H4'	23:DB:533:G:C8	2.35	0.62
23:DB:764:A:H5''	25:DC:208:GLY:HA2	1.82	0.62
47:DF:146:ASP:HB3	47:DF:147:ARG:NH1	2.14	0.62
40:DH:95:GLY:H	40:DH:98:ASP:HB2	1.64	0.62
41:DJ:29:ALA:HA	41:DJ:32:LEU:HD12	1.80	0.62
41:DJ:43:GLU:O	41:DJ:45:THR:N	2.32	0.62
38:DM:19:GLY:N	38:DM:38:ARG:HH12	1.97	0.62
52:DW:32:ALA:C	52:DW:34:SER:H	2.03	0.62
1:AA:1484:C:H2'	1:AA:1485:U:C6	2.35	0.61
1:AA:202:G:H21	1:AA:465:A:N6	1.97	0.61
1:AA:313:A:H2'	1:AA:314:C:C6	2.35	0.61
1:AA:474:G:H2'	1:AA:475:C:C6	2.35	0.61
1:AA:522:C:H41	11:AL:49:ARG:NH2	1.98	0.61
1:AA:736:C:H2'	1:AA:737:C:C6	2.34	0.61
1:AA:834:U:H2'	1:AA:835:U:C6	2.34	0.61
18:AB:132:GLU:O	18:AB:136:ARG:HG2	2.00	0.61
8:AI:82:ILE:O	8:AI:86:LEU:HD13	2.00	0.61
20:AO:69:TYR:HA	20:AO:72:ARG:NH1	2.14	0.61
1:AA:1319:A:H3'	16:AS:2:ARG:HA	1.82	0.61
16:AS:5:LYS:C	16:AS:6:LYS:HD2	2.20	0.61
19:AU:19:LYS:HD2	19:AU:19:LYS:N	2.15	0.61
23:BB:155:A:H2'	23:BB:156:A:H8	1.65	0.61
23:BB:1889:A:H2'	23:BB:1890:A:H8	1.65	0.61
25:BC:156:SER:O	25:BC:194:VAL:HG11	2.00	0.61
29:BE:158:PHE:O	29:BE:162:ARG:HB2	2.00	0.61
47:BF:48:LEU:H	47:BF:48:LEU:HD23	1.65	0.61
48:BG:120:ILE:HG13	48:BG:140:ILE:HG22	1.82	0.61
48:BG:15:ASP:HB2	48:BG:26:LYS:CD	2.29	0.61
48:BG:36:LEU:H	48:BG:36:LEU:HD22	1.63	0.61
24:BI:18:ASN:N	24:BI:19:PRO:HD2	2.14	0.61
27:BK:19:VAL:C	27:BK:41:ILE:HD11	2.20	0.61
37:BL:19:LEU:HD23	37:BL:27:LEU:HD23	1.82	0.61
37:BL:95:LEU:HB2	37:BL:101:ILE:CG1	2.30	0.61
27:BK:73:ASP:O	28:BP:74:GLN:HG3	2.00	0.61
35:BV:14:LYS:HE3	35:BV:18:ARG:HH21	1.65	0.61
1:CA:1320:C:O2'	1:CA:1321:U:H5'	2.00	0.61
1:CA:621:A:H2'	1:CA:622:A:C8	2.35	0.61
1:CA:704:A:C2	1:CA:705:G:H1'	2.35	0.61
1:CA:922:G:H2'	1:CA:923:A:C8	2.34	0.61
18:CB:216:VAL:HG23	18:CB:217:ALA:N	2.13	0.61
6:CG:130:LYS:H	6:CG:134:VAL:HG21	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CI:114:LYS:HE2	8:CI:120:ALA:HB1	1.82	0.61
10:CK:70:ALA:HA	10:CK:73:VAL:HG22	1.81	0.61
19:CU:13:VAL:HG22	19:CU:15:LEU:HD23	1.81	0.61
34:D3:14:LYS:HD3	34:D3:21:PHE:O	2.00	0.61
23:DB:1028:A:H2'	23:DB:1029:A:C8	2.35	0.61
23:DB:1279:G:H4'	42:DN:31:HIS:CD2	2.35	0.61
23:DB:1739:A:H2'	23:DB:1740:G:O4'	2.00	0.61
23:DB:2108:A:N3	23:DB:2108:A:H2'	2.13	0.61
23:DB:2786:U:H4'	26:DD:67:HIS:HA	1.82	0.61
23:DB:458:G:N2	23:DB:469:G:H2'	2.14	0.61
23:DB:544:C:HO2'	23:DB:546:U:H6	1.48	0.61
23:DB:968:C:H2'	23:DB:969:G:H8	1.65	0.61
48:DG:17:LYS:HE2	48:DG:19:ASN:ND2	2.14	0.61
48:DG:18:ILE:HA	48:DG:23:ILE:HG13	1.80	0.61
48:DG:21:GLN:OE1	48:DG:37:ASN:HB2	1.99	0.61
42:DN:87:PHE:HB3	42:DN:90:ARG:HB2	1.83	0.61
1:AA:1231:G:H2'	1:AA:1232:U:C6	2.35	0.61
1:AA:628:G:H2'	1:AA:629:A:C8	2.35	0.61
1:AA:955:U:H2'	1:AA:956:U:O4'	1.99	0.61
3:AD:197:HIS:O	3:AD:200:VAL:HG22	1.99	0.61
17:AT:4:LYS:HD2	17:AT:5:SER:N	2.15	0.61
23:BB:2860:A:O5'	23:BB:2860:A:H8	1.84	0.61
23:BB:845:A:C2	23:BB:847:U:H1'	2.35	0.61
41:BJ:58:ASN:HA	41:BJ:126:ALA:O	2.00	0.61
35:BV:80:HIS:CD2	35:BV:83:LYS:HB2	2.34	0.61
1:CA:1352:C:H2'	1:CA:1353:G:C8	2.35	0.61
1:CA:215:C:H2'	1:CA:216:U:C6	2.34	0.61
1:CA:22:G:H2'	1:CA:23:C:C6	2.35	0.61
1:CA:272:C:H2'	1:CA:273:U:H6	1.63	0.61
1:CA:810:C:O2'	1:CA:811:C:H5'	1.99	0.61
23:DB:1054:A:H2'	23:DB:1055:G:H8	1.63	0.61
23:DB:181:A:H2'	23:DB:182:A:C8	2.35	0.61
23:DB:2741:A:H2'	23:DB:2742:G:O4'	2.00	0.61
23:DB:351:C:H2'	23:DB:352:A:H8	1.65	0.61
25:DC:124:LYS:HE2	25:DC:127:ASN:HD21	1.64	0.61
41:DJ:58:ASN:HA	41:DJ:126:ALA:O	2.00	0.61
30:DY:23:LEU:HD13	30:DY:28:LEU:HB2	1.81	0.61
1:AA:1070:U:H2'	1:AA:1071:C:C6	2.36	0.61
1:AA:1366:C:H2'	1:AA:1367:C:H6	1.65	0.61
1:AA:1513:A:H2'	1:AA:1514:G:H8	1.63	0.61
1:AA:394:G:H2'	1:AA:395:C:C6	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:408:A:H3'	1:AA:409:U:H6	1.65	0.61
1:AA:764:C:C2'	1:AA:765:G:H5'	2.31	0.61
3:AD:146:GLU:HA	3:AD:149:LYS:HG2	1.81	0.61
6:AG:94:ARG:HH11	6:AG:98:LEU:HD11	1.64	0.61
23:BB:1778:U:H2'	23:BB:1784:A:H62	1.65	0.61
23:BB:2065:C:H2'	23:BB:2066:C:C6	2.35	0.61
23:BB:551:G:H2'	23:BB:552:U:C6	2.35	0.61
26:BD:55:LYS:NZ	26:BD:60:VAL:HG13	2.15	0.61
23:BB:615:U:O4	29:BE:39:ALA:HB2	2.00	0.61
24:BI:75:ALA:HB2	24:BI:112:LYS:HE2	1.81	0.61
38:BM:41:LEU:HB3	38:BM:46:ILE:CG2	2.30	0.61
50:BT:34:VAL:HG21	50:BT:43:ILE:HD12	1.83	0.61
51:BZ:35:SER:HA	51:BZ:50:ARG:HA	1.81	0.61
18:CB:56:LEU:HB3	18:CB:220:VAL:HG22	1.82	0.61
2:CC:82:ASP:O	2:CC:86:LEU:HG	2.01	0.61
8:CI:43:ALA:O	8:CI:47:VAL:HG23	1.99	0.61
11:CL:42:LYS:HD2	11:CL:43:LYS:HG2	1.82	0.61
17:CT:4:LYS:HD2	17:CT:5:SER:N	2.15	0.61
17:CT:8:LYS:HA	17:CT:11:ILE:HD12	1.82	0.61
22:DA:75:G:H1	22:DA:102:G:H22	1.48	0.61
23:DB:2104:C:H6	23:DB:2104:C:H3'	1.65	0.61
25:DC:115:ILE:HG22	25:DC:124:LYS:HE3	1.82	0.61
25:DC:92:LEU:HD12	25:DC:93:VAL:H	1.65	0.61
26:DD:116:LYS:HB3	26:DD:118:PHE:CZ	2.34	0.61
29:DE:158:PHE:O	29:DE:162:ARG:HB2	2.00	0.61
47:DF:62:GLN:CD	47:DF:91:ARG:HE	2.03	0.61
48:DG:145:ALA:O	48:DG:148:ARG:HG3	2.01	0.61
40:DH:96:THR:HG23	40:DH:97:ARG:H	1.66	0.61
24:DI:25:PRO:O	24:DI:29:GLN:HG3	2.00	0.61
27:DK:13:ASN:HD21	27:DK:98:ARG:HB2	1.65	0.61
37:DL:119:PRO:HA	37:DL:138:ALA:O	2.00	0.61
42:DN:94:TYR:O	42:DN:115:LEU:HA	2.00	0.61
1:AA:1243:C:H2'	1:AA:1244:G:H8	1.65	0.61
1:AA:1492:A:H3'	1:AA:1493:A:H4'	1.83	0.61
1:AA:478:A:H2'	1:AA:479:U:O4'	2.00	0.61
18:AB:80:LYS:O	18:AB:84:LEU:HB3	2.00	0.61
6:AG:23:ALA:O	6:AG:26:VAL:HG22	2.00	0.61
6:AG:30:MET:HG2	6:AG:31:VAL:H	1.65	0.61
21:AN:60:ARG:HD2	21:AN:69:PRO:HG3	1.81	0.61
34:B3:44:ARG:N	34:B3:45:PRO:HD2	2.15	0.61
23:BB:1774:C:O2	23:BB:1774:C:H2'	1.98	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1983:G:H4'	23:BB:2606:C:H4'	1.82	0.61
23:BB:2760:C:O2'	23:BB:2761:A:H5'	1.99	0.61
23:BB:615:U:O2	29:BE:35:TYR:HA	2.00	0.61
23:BB:82:U:H2'	23:BB:83:A:C8	2.35	0.61
23:BB:910:A:H2'	23:BB:911:A:H8	1.65	0.61
47:BF:22:ASN:ND2	47:BF:22:ASN:H	1.97	0.61
40:BH:68:ARG:O	40:BH:72:ILE:HG22	1.99	0.61
41:BJ:82:GLY:O	41:BJ:84:ILE:HG22	2.01	0.61
28:BP:13:LYS:HG2	28:BP:76:HIS:ND1	2.15	0.61
45:BS:15:GLN:HA	45:BS:18:ARG:CG	2.30	0.61
1:CA:1060:U:H5''	9:CJ:53:ILE:HG12	1.81	0.61
1:CA:157:U:O2'	1:CA:158:G:H5'	2.00	0.61
1:CA:602:A:O2'	1:CA:603:U:H5'	2.00	0.61
1:CA:677:U:H2'	1:CA:678:U:C6	2.35	0.61
18:CB:67:LEU:HA	18:CB:89:PHE:O	1.99	0.61
3:CD:29:THR:HB	3:CD:30:LYS:NZ	2.16	0.61
36:D2:31:LEU:HD22	36:D2:42:LEU:HD12	1.83	0.61
23:DB:1443:U:H2'	23:DB:1444:G:C8	2.36	0.61
23:DB:2547:A:H2'	23:DB:2548:U:C6	2.35	0.61
40:DH:121:VAL:HG23	40:DH:122:LEU:H	1.65	0.61
40:DH:96:THR:OG1	40:DH:112:LYS:HB2	2.00	0.61
42:DN:38:LEU:HD11	42:DN:42:LYS:HD2	1.83	0.61
46:DU:34:ILE:HG12	46:DU:63:ALA:HB2	1.81	0.61
1:AA:1138:G:H2'	1:AA:1139:G:H4'	1.82	0.61
1:AA:501:C:H2'	1:AA:502:A:H8	1.65	0.61
7:AH:44:PHE:HE2	7:AH:100:ILE:HG12	1.66	0.61
7:AH:86:LYS:HD2	7:AH:90:GLU:HG2	1.80	0.61
21:AN:60:ARG:CG	21:AN:62:ARG:HG2	2.28	0.61
19:AU:36:PHE:CB	19:AU:40:PRO:HD3	2.28	0.61
33:B1:3:GLY:O	33:B1:4:ILE:HG12	2.01	0.61
23:BB:1681:G:N3	23:BB:1762:A:H2'	2.15	0.61
23:BB:1870:C:H2'	23:BB:1871:A:C8	2.36	0.61
23:BB:2840:C:H2'	23:BB:2841:C:C6	2.35	0.61
23:BB:2861:U:H2'	23:BB:2862:G:H8	1.66	0.61
23:BB:2789:C:H3'	23:BB:2893:A:H62	1.63	0.61
47:BF:169:LEU:HA	47:BF:172:PHE:HD2	1.65	0.61
37:BL:121:THR:HB	37:BL:141:LYS:HB3	1.82	0.61
52:BW:24:ARG:HD3	52:BW:65:LYS:HE3	1.80	0.61
1:CA:1458:G:H2'	1:CA:1459:G:C8	2.34	0.61
18:CB:130:LYS:HA	18:CB:133:ALA:HB3	1.83	0.61
2:CC:26:LYS:CG	2:CC:27:GLU:H	2.10	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CH:58:LEU:HD22	7:CH:60:LEU:HB2	1.81	0.61
7:CH:74:ILE:HG13	7:CH:128:VAL:HG22	1.83	0.61
1:CA:522:C:H41	11:CL:49:ARG:NH2	1.97	0.61
16:CS:35:ARG:HB2	16:CS:71:GLY:HA2	1.82	0.61
10:CK:88:PRO:HD3	19:CU:28:LEU:HD11	1.83	0.61
23:DB:1180:U:H2'	23:DB:1181:U:C6	2.36	0.61
23:DB:2078:C:H2'	23:DB:2079:U:C6	2.36	0.61
23:DB:575:A:O2'	23:DB:576:U:H5'	2.00	0.61
23:DB:979:A:H2'	23:DB:982:C:N4	2.13	0.61
26:DD:55:LYS:NZ	26:DD:60:VAL:HG13	2.15	0.61
47:DF:84:ILE:HG23	47:DF:85:GLY:N	2.14	0.61
47:DF:94:ARG:HA	47:DF:97:GLU:HB3	1.82	0.61
38:DM:50:ARG:O	38:DM:53:MET:HB3	2.00	0.61
42:DN:33:ILE:HB	42:DN:114:GLU:HB2	1.83	0.61
49:DR:48:LYS:O	49:DR:48:LYS:HD2	2.01	0.61
1:AA:977:A:C2	1:AA:1223:C:H2'	2.36	0.61
2:AC:13:ILE:O	2:AC:14:VAL:HG22	2.00	0.61
6:AG:44:SER:OG	6:AG:116:ALA:HB1	2.01	0.61
6:AG:14:ASP:HB3	6:AG:18:GLY:H	1.64	0.61
14:AQ:46:HIS:HB2	14:AQ:70:LYS:HE2	1.81	0.61
36:B2:27:GLY:O	36:B2:30:VAL:HB	2.00	0.61
23:BB:152:A:H2'	23:BB:153:U:C6	2.35	0.61
23:BB:1709:U:H2'	23:BB:1710:G:C8	2.35	0.61
23:BB:172:A:O2'	23:BB:173:A:H5'	2.01	0.61
23:BB:2626:C:H2'	23:BB:2627:G:C8	2.36	0.61
25:BC:115:ILE:HG22	25:BC:124:LYS:HE3	1.83	0.61
47:BF:32:LYS:HA	47:BF:95:MET:HG3	1.82	0.61
40:BH:78:VAL:O	40:BH:143:ILE:HG13	1.98	0.61
24:BI:20:SER:O	24:BI:25:PRO:HD2	2.01	0.61
27:BK:25:LEU:HD21	27:BK:40:LYS:H	1.66	0.61
28:BP:20:ARG:HG3	28:BP:21:PRO:CD	2.31	0.61
44:BQ:91:ARG:HD3	49:BR:11:GLN:HB2	1.81	0.61
46:BU:85:ARG:HD3	46:BU:86:PHE:H	1.66	0.61
30:BY:15:ARG:O	30:BY:20:LYS:HE3	2.00	0.61
1:CA:1448:C:H2'	1:CA:1449:C:C6	2.36	0.61
1:CA:543:U:H2'	1:CA:544:G:C8	2.36	0.61
1:CA:593:U:H2'	1:CA:594:U:C6	2.36	0.61
9:CJ:43:PRO:O	9:CJ:71:LEU:HD21	2.00	0.61
16:CS:4:LEU:HD13	16:CS:8:PRO:HA	1.81	0.61
1:CA:1221:G:C4'	16:CS:76:THR:HG21	2.28	0.61
23:DB:1889:A:H2'	23:DB:1890:A:H8	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DE:58:LYS:HB2	29:DE:60:TRP:CD1	2.35	0.61
47:DF:137:PHE:O	47:DF:139:GLU:N	2.33	0.61
48:DG:120:ILE:HG13	48:DG:140:ILE:HG22	1.82	0.61
48:DG:15:ASP:HB3	48:DG:26:LYS:H	1.66	0.61
48:DG:84:LYS:HG3	48:DG:131:VAL:C	2.20	0.61
23:DB:1142:A:H4'	41:DJ:27:ARG:HH22	1.64	0.61
23:DB:2379:G:H4'	43:DO:21:LEU:HD11	1.81	0.61
45:DS:20:VAL:C	45:DS:22:ASP:H	2.03	0.61
46:DU:11:ILE:O	46:DU:11:ILE:HD13	2.01	0.61
2:AC:96:VAL:HB	2:AC:97:PRO:HD2	1.81	0.61
11:AL:42:LYS:HD2	11:AL:43:LYS:HG2	1.83	0.61
23:BB:1203:U:H3'	23:BB:1204:A:C5'	2.31	0.61
23:BB:1534:U:H2'	23:BB:1536:C:N3	2.15	0.61
23:BB:2678:C:H2'	23:BB:2679:A:C8	2.35	0.61
23:BB:6:A:H2'	23:BB:7:G:H8	1.65	0.61
25:BC:6:LYS:O	25:BC:8:THR:HG23	2.00	0.61
47:BF:146:ASP:HB3	47:BF:147:ARG:NH1	2.15	0.61
40:BH:6:LEU:H	40:BH:6:LEU:HD12	1.65	0.61
40:BH:83:LYS:HB3	40:BH:91:PHE:CD1	2.35	0.61
41:BJ:4:PHE:HB3	41:BJ:44:TYR:CD1	2.35	0.61
35:BV:62:THR:HG22	35:BV:71:LYS:HG2	1.82	0.61
52:BW:18:LYS:HG3	52:BW:19:ARG:CZ	2.31	0.61
1:CA:1132:C:H2'	1:CA:1133:G:C8	2.35	0.61
11:CL:35:ARG:NH2	11:CL:36:VAL:HG22	2.16	0.61
20:CO:32:LEU:O	20:CO:36:ILE:HG12	2.01	0.61
23:DB:2046:G:H1'	31:D0:18:HIS:CD2	2.34	0.61
34:D3:61:LEU:N	34:D3:62:PRO:HD3	2.15	0.61
23:DB:1203:U:H1'	37:DL:4:ASN:HD21	1.64	0.61
23:DB:18:U:H2'	23:DB:19:A:H8	1.66	0.61
23:DB:2393:U:H5''	37:DL:62:PRO:HG3	1.82	0.61
23:DB:2737:G:H2'	23:DB:2738:A:C8	2.35	0.61
26:DD:119:ALA:HB1	26:DD:163:GLY:H	1.66	0.61
47:DF:78:ILE:HA	47:DF:82:TYR:CG	2.36	0.61
23:DB:2376:A:N1	43:DO:92:PHE:HD2	1.99	0.61
1:AA:1508:A:H2'	1:AA:1509:C:C6	2.36	0.61
1:AA:624:C:H2'	1:AA:625:U:H6	1.64	0.61
2:AC:137:VAL:HG22	2:AC:150:VAL:HG21	1.83	0.61
32:B4:1:MET:HE1	32:B4:36:ARG:HB2	1.82	0.61
23:BB:1198:U:H2'	23:BB:1199:U:C6	2.35	0.61
23:BB:1821:A:H2'	23:BB:1822:C:H6	1.64	0.61
23:BB:2352:A:H2'	23:BB:2353:G:O4'	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:401:A:H2'	23:BB:402:A:C8	2.36	0.61
23:BB:5:A:H2'	23:BB:6:A:C8	2.35	0.61
26:BD:54:ALA:CA	26:BD:76:GLY:HA2	2.31	0.61
47:BF:62:GLN:CD	47:BF:91:ARG:HE	2.03	0.61
24:BI:7:TYR:HB2	24:BI:58:ILE:O	2.00	0.61
27:BK:59:LYS:HD2	27:BK:89:ASN:HD22	1.65	0.61
37:BL:80:SER:H	37:BL:113:ALA:HB3	1.65	0.61
23:BB:811:U:H2'	37:BL:21:ARG:HA	1.81	0.61
1:CA:1347:G:H22	1:CA:1373:G:H2'	1.66	0.61
1:CA:209:U:H5'	1:CA:210:C:H5	1.65	0.61
1:CA:764:C:H2'	1:CA:765:G:H5'	1.82	0.61
1:CA:806:C:H2'	1:CA:807:A:H8	1.66	0.61
18:CB:86:CYS:H	18:CB:88:GLN:NE2	1.99	0.61
9:CJ:6:ILE:O	9:CJ:75:ASP:HA	2.00	0.61
23:DB:1105:U:H2'	23:DB:1106:G:C8	2.36	0.61
23:DB:1220:G:H2'	23:DB:1221:C:C6	2.36	0.61
38:DM:35:ALA:H	38:DM:100:LYS:H	1.48	0.61
45:DS:18:ARG:HB3	45:DS:76:VAL:HG22	1.83	0.61
23:DB:851:C:H4'	30:DY:46:MET:HA	1.83	0.61
51:DZ:4:VAL:O	51:DZ:33:LEU:HD23	2.00	0.61
51:DZ:67:VAL:O	51:DZ:70:GLU:HG2	2.01	0.61
1:AA:1316:G:N2	1:AA:1318:A:H3'	2.15	0.61
1:AA:593:U:H2'	1:AA:594:U:C6	2.36	0.61
1:AA:662:U:H2'	1:AA:663:A:H8	1.64	0.61
1:AA:764:C:H2'	1:AA:765:G:H5'	1.82	0.61
1:AA:781:A:H2'	1:AA:782:A:H5'	1.83	0.61
18:AB:14:HIS:HD2	18:AB:202:ASN:HB2	1.66	0.61
3:AD:121:ALA:O	3:AD:122:ILE:HD13	2.00	0.61
1:AA:522:C:H41	11:AL:49:ARG:HH22	1.48	0.61
13:AP:67:ILE:HG13	13:AP:71:VAL:HG13	1.83	0.61
17:AT:8:LYS:HA	17:AT:11:ILE:HD12	1.83	0.61
23:BB:126:A:O5'	36:B2:19:ARG:HG3	2.01	0.61
32:B4:16:ILE:HG12	32:B4:25:VAL:HG22	1.83	0.61
32:B4:7:VAL:HG13	32:B4:8:LYS:N	2.13	0.61
23:BB:1788:C:O2'	23:BB:1789:A:H5'	2.01	0.61
23:BB:635:C:O2'	23:BB:639:U:H5''	2.01	0.61
25:BC:77:VAL:HG23	25:BC:112:GLY:H	1.66	0.61
29:BE:67:ARG:HG2	29:BE:67:ARG:HH11	1.65	0.61
47:BF:34:THR:O	47:BF:89:THR:HA	2.01	0.61
27:BK:109:SER:HB2	27:BK:111:LYS:HE2	1.82	0.61
27:BK:70:ARG:HD3	27:BK:76:VAL:HG22	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BQ:80:ASN:O	44:BQ:83:LYS:HB3	2.01	0.61
50:BT:32:LEU:N	50:BT:83:ALA:HB3	2.16	0.61
46:BU:11:ILE:O	46:BU:11:ILE:HD13	2.00	0.61
1:CA:1000:A:H2'	1:CA:1001:C:C6	2.35	0.61
1:CA:1061:G:H2'	1:CA:1062:U:C6	2.36	0.61
1:CA:1492:A:H5''	1:CA:1493:A:C8	2.36	0.61
1:CA:26:A:H61	1:CA:558:G:H1'	1.65	0.61
18:CB:209:VAL:HG23	18:CB:210:THR:H	1.66	0.61
9:CJ:7:ARG:HE	9:CJ:101:SER:HB3	1.65	0.61
21:CN:62:ARG:HG2	21:CN:67:GLY:O	2.01	0.61
34:D3:44:ARG:N	34:D3:45:PRO:HD2	2.15	0.61
23:DB:1470:A:H3'	23:DB:1471:G:H8	1.66	0.61
23:DB:2573:C:H3'	56:DB:3628:HOH:O	2.01	0.61
26:DD:68:PHE:HB3	26:DD:73:VAL:HG23	1.81	0.61
24:DI:42:ASN:HA	24:DI:45:THR:OG1	1.99	0.61
37:DL:57:LEU:HD12	37:DL:60:ARG:NH2	2.16	0.61
1:AA:1308:U:H2'	1:AA:1309:G:H8	1.66	0.61
1:AA:312:C:H2'	1:AA:313:A:C8	2.35	0.61
14:AQ:16:MET:HB2	14:AQ:19:SER:HB2	1.81	0.61
23:BB:2073:C:O2'	23:BB:2074:U:H5'	2.00	0.61
23:BB:2741:A:H2'	23:BB:2742:G:O4'	2.01	0.61
23:BB:2747:G:O6	23:BB:2754:U:H2'	2.01	0.61
26:BD:8:LYS:HG2	26:BD:197:THR:H	1.66	0.61
47:BF:177:ARG:NH1	47:BF:177:ARG:HA	2.15	0.61
40:BH:44:ILE:HA	40:BH:51:ARG:HH22	1.66	0.61
44:BQ:100:PHE:HB2	49:BR:13:ARG:HH12	1.65	0.61
45:BS:5:ALA:HB3	45:BS:54:ALA:HB2	1.83	0.61
1:CA:1222:G:O2'	1:CA:1223:C:H5'	2.01	0.61
1:CA:1308:U:O2'	1:CA:1309:G:H5'	2.01	0.61
1:CA:1342:C:H2'	1:CA:1343:G:C8	2.36	0.61
1:CA:373:A:H2'	1:CA:374:A:H8	1.66	0.61
1:CA:202:G:H21	1:CA:465:A:N6	1.99	0.61
23:DB:1054:A:H2'	23:DB:1055:G:C8	2.36	0.61
23:DB:2352:A:H2'	23:DB:2353:G:O4'	2.01	0.61
47:DF:134:GLN:C	47:DF:136:ILE:H	2.05	0.61
24:DI:20:SER:O	24:DI:25:PRO:HD2	2.00	0.61
27:DK:85:VAL:HG21	27:DK:115:ILE:HD11	1.82	0.61
38:DM:18:ARG:C	38:DM:38:ARG:HH22	2.03	0.61
23:DB:1599:U:OP1	50:DT:40:LYS:HB3	2.01	0.61
1:AA:977:A:H2	1:AA:1223:C:H2'	1.66	0.60
1:AA:464:U:H3'	1:AA:466:A:OP1	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:946:A:H2'	1:AA:947:G:H8	1.65	0.60
4:AE:131:ASN:HD21	4:AE:133:ILE:HB	1.64	0.60
6:AG:138:GLU:HB3	6:AG:142:ARG:NH2	2.15	0.60
21:AN:11:LYS:O	21:AN:15:LEU:HG	2.00	0.60
20:AO:67:LEU:HD12	20:AO:88:ARG:HH22	1.65	0.60
33:B1:36:LYS:HB2	33:B1:47:ILE:HA	1.83	0.60
23:BB:1220:G:H2'	23:BB:1221:C:H6	1.65	0.60
23:BB:1443:U:H2'	23:BB:1444:G:C8	2.35	0.60
23:BB:1739:A:H2'	23:BB:1740:G:O4'	2.00	0.60
23:BB:2216:G:H2'	23:BB:2217:G:C8	2.35	0.60
23:BB:2742:G:OP1	32:B4:36:ARG:HD3	2.01	0.60
23:BB:834:G:H2'	23:BB:835:C:O4'	2.01	0.60
25:BC:94:LEU:HA	25:BC:100:ARG:HA	1.83	0.60
47:BF:111:ARG:HE	47:BF:135:ILE:HG22	1.66	0.60
24:BI:11:GLN:HA	24:BI:55:PRO:HA	1.82	0.60
37:BL:119:PRO:HA	37:BL:138:ALA:O	2.01	0.60
35:BV:80:HIS:CD2	35:BV:83:LYS:H	2.19	0.60
1:CA:1226:C:O2'	1:CA:1227:A:H5'	2.01	0.60
1:CA:478:A:H2'	1:CA:479:U:O4'	2.01	0.60
5:CF:71:ILE:HG13	5:CF:72:ASP:N	2.14	0.60
11:CL:41:PRO:HB3	11:CL:49:ARG:NH1	2.16	0.60
12:CM:102:LYS:HG3	12:CM:103:THR:HG23	1.83	0.60
14:CQ:16:MET:HB2	14:CQ:19:SER:HB2	1.82	0.60
23:DB:1175:A:C4	23:DB:1176:U:H1'	2.36	0.60
23:DB:1406:U:H2'	23:DB:1407:G:C8	2.35	0.60
23:DB:1782:U:H3'	56:DB:3606:HOH:O	2.01	0.60
23:DB:263:G:H2'	23:DB:264:C:O4'	2.01	0.60
23:DB:2677:G:H2'	23:DB:2678:C:C6	2.35	0.60
23:DB:277:G:H1'	23:DB:361:G:O6	2.01	0.60
23:DB:2840:C:H2'	23:DB:2841:C:C6	2.34	0.60
26:DD:30:GLU:HB2	26:DD:52:THR:HG23	1.83	0.60
47:DF:120:SER:HG	47:DF:127:TYR:HD2	1.48	0.60
46:DU:42:LYS:HG3	46:DU:57:ILE:HG21	1.83	0.60
1:AA:462:G:H3'	1:AA:463:U:H6	1.66	0.60
1:AA:675:A:H2'	1:AA:676:A:C8	2.36	0.60
1:AA:797:C:O2'	1:AA:798:U:H5'	2.00	0.60
18:AB:130:LYS:O	18:AB:134:LEU:HG	2.01	0.60
2:AC:8:GLY:HA2	2:AC:11:LEU:HD21	1.81	0.60
8:AI:96:GLU:HA	8:AI:99:LYS:NZ	2.16	0.60
22:BA:6:G:H2'	22:BA:7:G:H8	1.64	0.60
23:BB:1631:G:N2	23:BB:1633:G:H3'	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1868:C:H2'	23:BB:1869:G:O4'	2.01	0.60
23:BB:2516:A:O2'	23:BB:2517:C:H5'	2.01	0.60
23:BB:263:G:H2'	23:BB:264:C:O4'	2.01	0.60
23:BB:2788:C:H2'	23:BB:2789:C:C6	2.35	0.60
23:BB:918:A:H2'	23:BB:919:U:H5'	1.82	0.60
26:BD:10:GLY:O	26:BD:11:MET:HB2	2.00	0.60
26:BD:39:ASP:HB3	26:BD:42:ASN:HB3	1.83	0.60
29:BE:145:ASP:HB3	29:BE:184:ASP:H	1.67	0.60
40:BH:117:LEU:HD13	40:BH:120:GLY:HA2	1.83	0.60
41:BJ:124:VAL:HG23	41:BJ:125:TYR:H	1.64	0.60
41:BJ:96:ARG:O	41:BJ:99:ARG:HG3	2.01	0.60
35:BV:80:HIS:HD2	35:BV:83:LYS:H	1.46	0.60
52:BW:44:PHE:O	52:BW:78:PHE:HA	2.00	0.60
1:CA:1316:G:H22	1:CA:1318:A:H3'	1.65	0.60
1:CA:188:C:H2'	1:CA:189:A:O4'	2.01	0.60
1:CA:279:A:H5'	1:CA:281:G:H5'	1.82	0.60
1:CA:860:A:H2'	1:CA:861:G:O4'	2.01	0.60
2:CC:59:PRO:CG	2:CC:62:SER:HB2	2.28	0.60
21:CN:16:ALA:HB2	21:CN:54:SER:O	2.00	0.60
32:D4:22:VAL:O	32:D4:24:ARG:HG3	2.00	0.60
23:DB:2078:C:O2'	23:DB:2079:U:H5'	2.01	0.60
23:DB:2569:G:O2'	23:DB:2570:G:H5'	2.01	0.60
23:DB:438:G:H2'	23:DB:439:A:H8	1.66	0.60
23:DB:971:G:H2'	23:DB:972:A:O4'	2.00	0.60
25:DC:130:PRO:HA	25:DC:188:ARG:HA	1.83	0.60
24:DI:78:LEU:HA	24:DI:81:LYS:HE2	1.82	0.60
37:DL:2:ARG:HG2	37:DL:5:THR:HG21	1.82	0.60
35:DV:75:GLN:HG2	35:DV:92:VAL:HG23	1.82	0.60
23:DB:2264:C:H41	52:DW:11:ASN:ND2	1.99	0.60
1:AA:1047:G:N2	1:AA:1215:G:H5''	2.16	0.60
1:AA:1053:G:O6	1:AA:1199:U:H2'	2.01	0.60
1:AA:1163:A:H2'	1:AA:1164:G:C8	2.35	0.60
1:AA:462:G:H3'	1:AA:463:U:C6	2.36	0.60
1:AA:923:A:H2'	1:AA:924:C:C6	2.36	0.60
18:AB:35:ASN:C	18:AB:36:LYS:HD2	2.21	0.60
7:AH:55:LYS:HA	7:AH:55:LYS:HE3	1.83	0.60
20:AO:39:LEU:HD23	20:AO:56:LEU:HB2	1.82	0.60
23:BB:1082:U:N3	23:BB:1086:A:C6	2.70	0.60
23:BB:1315:C:H2'	23:BB:1316:U:C6	2.36	0.60
23:BB:1485:U:H2'	23:BB:1486:U:H6	1.66	0.60
23:BB:2103:C:H2'	23:BB:2104:C:O4'	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:480:A:H3'	23:BB:481:G:H5'	1.83	0.60
23:BB:633:A:O5'	23:BB:633:A:H8	1.84	0.60
23:BB:850:U:H2'	23:BB:851:C:C6	2.36	0.60
24:BI:89:SER:HA	24:BI:97:VAL:HG21	1.83	0.60
23:BB:558:U:OP1	41:BJ:114:LEU:HB2	2.01	0.60
37:BL:68:SER:HB2	37:BL:71:ALA:HB3	1.83	0.60
45:BS:20:VAL:C	45:BS:22:ASP:H	2.05	0.60
1:CA:441:A:H61	1:CA:493:A:N6	2.00	0.60
1:CA:806:C:H2'	1:CA:807:A:C8	2.37	0.60
18:CB:46:VAL:CG1	18:CB:47:PRO:HD3	2.26	0.60
3:CD:169:TRP:O	3:CD:182:LYS:HB2	2.01	0.60
7:CH:94:VAL:HG12	7:CH:99:GLY:HA3	1.83	0.60
23:DB:1098:A:C8	24:DI:3:LYS:HB3	2.36	0.60
23:DB:155:A:H2'	23:DB:156:A:H8	1.65	0.60
23:DB:2065:C:H2'	23:DB:2066:C:C6	2.36	0.60
23:DB:401:A:H2'	23:DB:402:A:C8	2.35	0.60
25:DC:90:ILE:HD13	25:DC:104:LEU:HA	1.83	0.60
29:DE:60:TRP:CH2	29:DE:69:ARG:HA	2.36	0.60
24:DI:27:LEU:CD2	24:DI:27:LEU:H	2.14	0.60
45:DS:15:GLN:HA	45:DS:18:ARG:CG	2.32	0.60
1:AA:840:C:C4	1:AA:842:U:H4'	2.37	0.60
2:AC:87:ARG:NH2	2:AC:88:LYS:HA	2.16	0.60
3:AD:29:THR:H	3:AD:33:ILE:CG2	2.13	0.60
14:AQ:18:LYS:H	14:AQ:50:ASN:HD21	1.47	0.60
36:B2:20:ALA:C	36:B2:22:MET:H	2.05	0.60
23:BB:1577:C:H2'	23:BB:1578:U:C6	2.37	0.60
23:BB:2261:C:H1'	23:BB:2388:A:N3	2.16	0.60
23:BB:654:A:N3	23:BB:654:A:H2'	2.15	0.60
47:BF:177:ARG:CZ	47:BF:177:ARG:HA	2.32	0.60
27:BK:38:ILE:H	27:BK:38:ILE:CD1	2.14	0.60
49:BR:36:ALA:HB2	49:BR:58:VAL:HG23	1.83	0.60
49:BR:49:ILE:HD12	49:BR:49:ILE:O	2.02	0.60
46:BU:26:ASN:HD22	46:BU:26:ASN:N	1.98	0.60
51:BZ:67:VAL:O	51:BZ:70:GLU:HG2	2.02	0.60
18:CB:172:ILE:HG22	18:CB:176:ASN:ND2	2.17	0.60
18:CB:60:ALA:HB1	18:CB:224:ARG:HG2	1.82	0.60
6:CG:105:GLU:HB3	6:CG:136:LYS:NZ	2.16	0.60
6:CG:139:ASP:HA	6:CG:142:ARG:HH11	1.65	0.60
21:CN:27:LYS:CG	21:CN:28:ALA:H	2.13	0.60
21:CN:63:CYS:SG	21:CN:66:THR:N	2.74	0.60
16:CS:35:ARG:HA	16:CS:70:LEU:HB3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1199:U:H2'	23:DB:1200:C:H6	1.66	0.60
23:DB:1551:A:H3'	23:DB:1552:A:H5''	1.81	0.60
23:DB:2469:A:H4'	38:DM:55:ARG:CZ	2.32	0.60
23:DB:2698:U:H2'	23:DB:2699:C:C6	2.35	0.60
23:DB:2824:C:H3'	23:DB:2825:G:N2	2.17	0.60
23:DB:6:A:H4'	41:DJ:131:ASN:O	2.01	0.60
23:DB:848:C:H2'	23:DB:849:A:C8	2.35	0.60
26:DD:202:ILE:HG22	26:DD:204:LYS:NZ	2.16	0.60
44:DQ:87:VAL:HG21	49:DR:52:PRO:HG3	1.81	0.60
45:DS:66:ILE:N	45:DS:66:ILE:HD13	2.16	0.60
50:DT:14:PRO:HA	50:DT:32:LEU:CB	2.32	0.60
50:DT:57:VAL:HG22	50:DT:58:VAL:N	2.14	0.60
35:DV:80:HIS:HD2	35:DV:83:LYS:H	1.49	0.60
1:AA:602:A:O2'	1:AA:603:U:H5'	2.02	0.60
1:AA:621:A:H2'	1:AA:622:A:C8	2.36	0.60
6:AG:45:ALA:HB1	6:AG:120:ALA:HA	1.83	0.60
8:AI:117:LEU:HD12	8:AI:120:ALA:O	2.01	0.60
32:B4:27:CYS:SG	32:B4:29:ALA:HB3	2.41	0.60
23:BB:1117:C:H2'	23:BB:1118:C:C6	2.36	0.60
23:BB:1319:C:O2'	23:BB:1320:C:H5'	2.01	0.60
23:BB:2267:A:H8	23:BB:2267:A:H3'	1.66	0.60
23:BB:2292:U:H2'	23:BB:2293:G:H8	1.67	0.60
23:BB:458:G:N2	23:BB:469:G:H2'	2.16	0.60
23:BB:968:C:H2'	23:BB:969:G:H8	1.65	0.60
41:BJ:58:ASN:N	41:BJ:127:GLY:HA2	2.16	0.60
1:CA:32:A:H2'	1:CA:33:A:C8	2.36	0.60
1:CA:784:A:H2'	1:CA:785:G:H8	1.67	0.60
9:CJ:29:ALA:O	9:CJ:32:THR:HG22	2.01	0.60
23:DB:1593:A:H2'	23:DB:1594:U:C6	2.37	0.60
23:DB:2136:G:O2'	23:DB:2137:U:H5'	2.01	0.60
23:DB:2272:U:HO2'	23:DB:2273:A:H8	1.48	0.60
23:DB:2341:G:H2'	23:DB:2342:C:H6	1.66	0.60
23:DB:264:C:C2'	23:DB:265:A:H5''	2.32	0.60
23:DB:477:A:H2'	23:DB:478:A:H8	1.63	0.60
23:DB:620:G:N3	23:DB:620:G:H5'	2.16	0.60
23:DB:870:U:H2'	23:DB:871:U:H5''	1.83	0.60
23:DB:912:C:O2'	23:DB:913:U:H5'	2.01	0.60
26:DD:108:ASP:OD2	26:DD:173:GLN:HA	2.02	0.60
47:DF:77:LYS:HD2	47:DF:79:ARG:HD3	1.82	0.60
23:DB:1063:G:H4'	24:DI:135:MET:HG2	1.83	0.60
41:DJ:82:GLY:O	41:DJ:84:ILE:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DQ:90:ASP:OD1	49:DR:39:LEU:HD12	2.00	0.60
39:DX:12:GLU:HA	39:DX:15:ASN:HD21	1.66	0.60
1:AA:1148:U:H2'	1:AA:1149:C:H5'	1.83	0.60
18:AB:162:VAL:HG21	18:AB:168:GLU:HB2	1.82	0.60
2:AC:141:MET:HA	2:AC:141:MET:HE2	1.82	0.60
2:AC:184:ASN:HD22	2:AC:185:THR:H	1.49	0.60
2:AC:95:GLY:O	2:AC:96:VAL:HG13	2.00	0.60
3:AD:169:TRP:O	3:AD:182:LYS:HB2	2.02	0.60
3:AD:29:THR:HG22	3:AD:30:LYS:N	2.16	0.60
5:AF:53:LYS:HB2	5:AF:54:LEU:HD22	1.82	0.60
21:AN:58:ARG:H	21:AN:59:GLN:NE2	2.00	0.60
36:B2:16:HIS:HB3	36:B2:21:ARG:NH1	2.17	0.60
23:BB:125:A:H5'	36:B2:19:ARG:HD3	1.82	0.60
23:BB:1794:A:H2'	23:BB:1795:C:H6	1.66	0.60
23:BB:2041:U:H2'	23:BB:2042:A:C8	2.37	0.60
23:BB:2747:G:H2'	23:BB:2748:A:C8	2.37	0.60
23:BB:2859:G:H2'	23:BB:2860:A:C8	2.36	0.60
23:BB:455:C:N3	23:BB:472:A:H2'	2.16	0.60
23:BB:648:G:H2'	23:BB:649:G:H8	1.66	0.60
26:BD:2:ILE:HG22	26:BD:82:PHE:HB3	1.81	0.60
47:BF:62:GLN:NE2	47:BF:90:LEU:HD13	2.15	0.60
42:BN:48:VAL:HA	42:BN:51:LEU:HD22	1.84	0.60
44:BQ:84:LYS:HE2	44:BQ:84:LYS:HA	1.81	0.60
50:BT:43:ILE:O	50:BT:47:VAL:HG23	2.01	0.60
46:BU:51:LEU:H	46:BU:53:GLN:NE2	1.99	0.60
46:BU:85:ARG:NH1	46:BU:86:PHE:H	2.00	0.60
51:BZ:4:VAL:O	51:BZ:33:LEU:HD23	2.02	0.60
1:CA:135:C:C2	13:CP:1:MET:HB2	2.36	0.60
1:CA:279:A:C5'	1:CA:280:C:H3'	2.30	0.60
1:CA:628:G:H2'	1:CA:629:A:C8	2.36	0.60
1:CA:753:A:OP1	20:CO:73:LYS:HE3	2.01	0.60
2:CC:106:ARG:CZ	2:CC:106:ARG:H	2.14	0.60
6:CG:132:THR:HA	6:CG:135:LYS:HB3	1.84	0.60
16:CS:27:LYS:NZ	16:CS:27:LYS:HB3	2.15	0.60
22:DA:55:U:H2'	22:DA:56:G:C8	2.37	0.60
23:DB:1082:U:N3	23:DB:1086:A:C6	2.69	0.60
23:DB:1722:A:H2'	23:DB:1723:G:H8	1.66	0.60
23:DB:2262:U:H2'	23:DB:2263:C:C6	2.36	0.60
23:DB:2860:A:O5'	23:DB:2860:A:H8	1.82	0.60
23:DB:759:G:H2'	23:DB:760:G:H8	1.67	0.60
26:DD:104:VAL:HG12	26:DD:106:LYS:HE2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:168:GLU:O	26:DD:170:VAL:HG22	2.02	0.60
47:DF:134:GLN:O	47:DF:136:ILE:N	2.35	0.60
47:DF:62:GLN:NE2	47:DF:90:LEU:HD13	2.17	0.60
48:DG:157:LYS:HG2	48:DG:159:LYS:HG3	1.82	0.60
42:DN:85:PRO:HA	42:DN:88:ALA:HB2	1.84	0.60
50:DT:4:GLU:OE2	50:DT:5:GLU:HG2	2.01	0.60
39:DX:14:LEU:HD22	39:DX:57:LEU:HD21	1.84	0.60
1:AA:1262:C:N4	1:AA:1273:C:H42	1.98	0.60
1:AA:209:U:H5'	1:AA:210:C:C5	2.36	0.60
18:AB:107:ARG:HH21	18:AB:111:LYS:HB2	1.67	0.60
23:BB:1061:U:H4'	23:BB:1070:A:O3'	2.02	0.60
23:BB:1142:A:H4'	41:BJ:27:ARG:HH22	1.66	0.60
23:BB:1229:C:H2'	23:BB:1230:A:H8	1.66	0.60
23:BB:1810:A:H2'	23:BB:1811:G:O4'	2.01	0.60
23:BB:2262:U:H2'	23:BB:2263:C:C6	2.35	0.60
23:BB:479:A:N3	23:BB:481:G:H5''	2.17	0.60
26:BD:30:GLU:HB2	26:BD:52:THR:HG23	1.82	0.60
47:BF:24:VAL:O	47:BF:27:VAL:HG22	2.01	0.60
48:BG:137:LYS:O	48:BG:140:ILE:HG13	2.02	0.60
40:BH:78:VAL:HG12	40:BH:144:VAL:HG23	1.84	0.60
44:BQ:10:ARG:CZ	44:BQ:10:ARG:HB2	2.31	0.60
52:BW:49:ASN:HB3	52:BW:60:ALA:HA	1.84	0.60
1:CA:1492:A:H5'	1:CA:1493:A:OP2	2.02	0.60
1:CA:235:C:H2'	1:CA:236:A:C8	2.37	0.60
1:CA:436:C:O2'	1:CA:437:U:H5'	2.02	0.60
1:CA:553:A:H2'	1:CA:554:A:C8	2.37	0.60
1:CA:713:G:H2'	1:CA:714:G:H8	1.65	0.60
6:CG:38:ALA:O	6:CG:42:VAL:HG23	2.00	0.60
19:CU:20:ARG:HD2	19:CU:20:ARG:H	1.66	0.60
31:D0:53:VAL:HG12	42:DN:118:ARG:HH12	1.67	0.60
33:D1:32:LYS:HA	33:D1:51:ALA:O	2.01	0.60
23:DB:104:A:H2'	23:DB:105:C:C6	2.36	0.60
23:DB:2336:A:N6	52:DW:40:ARG:HG3	2.17	0.60
23:DB:2720:U:H5''	28:DP:52:ARG:NH2	2.15	0.60
23:DB:558:U:P	41:DJ:113:PRO:HG2	2.41	0.60
26:DD:4:LEU:HD23	26:DD:101:PHE:CE1	2.37	0.60
40:DH:6:LEU:HD12	40:DH:6:LEU:H	1.66	0.60
27:DK:43:ILE:HD12	27:DK:56:ASP:HB2	1.84	0.60
23:DB:2394:C:OP1	37:DL:63:LYS:HG2	2.01	0.60
49:DR:36:ALA:HB2	49:DR:58:VAL:HG23	1.84	0.60
46:DU:54:PRO:HG2	46:DU:55:GLY:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DX:13:GLU:O	39:DX:17:GLU:HG3	2.00	0.60
51:DZ:3:ARG:HG2	51:DZ:33:LEU:HD22	1.83	0.60
1:AA:1147:C:H2'	1:AA:1148:U:C6	2.37	0.60
1:AA:182:A:HO2'	1:AA:183:C:H3'	1.66	0.60
1:AA:215:C:H2'	1:AA:216:U:H6	1.66	0.60
18:AB:109:SER:O	18:AB:112:ARG:HB3	2.02	0.60
7:AH:74:ILE:HG13	7:AH:128:VAL:HG22	1.82	0.60
11:AL:17:LYS:N	11:AL:17:LYS:HE3	2.16	0.60
12:AM:72:ILE:O	12:AM:76:ILE:HG13	2.01	0.60
33:B1:5:ARG:NH1	33:B1:25:ASN:HB2	2.17	0.60
23:BB:2247:A:H3'	56:BB:3268:HOH:O	2.02	0.60
23:BB:2751:G:O2'	23:BB:2752:C:H5'	2.02	0.60
23:BB:62:U:H2'	23:BB:63:A:O4'	2.02	0.60
25:BC:92:LEU:HD12	25:BC:93:VAL:N	2.16	0.60
42:BN:33:ILE:HA	42:BN:114:GLU:HB2	1.83	0.60
44:BQ:4:LYS:HZ1	44:BQ:7:VAL:HG22	1.67	0.60
1:CA:1237:C:H3'	1:CA:1336:C:H41	1.66	0.60
1:CA:1292:G:H2'	1:CA:1293:C:H6	1.65	0.60
1:CA:1318:A:H4'	16:CS:9:PHE:CD1	2.36	0.60
1:CA:1350:A:H2'	1:CA:1351:U:C6	2.37	0.60
1:CA:313:A:H2'	1:CA:314:C:C6	2.37	0.60
1:CA:464:U:H3'	1:CA:466:A:OP1	2.02	0.60
1:CA:865:A:C2	1:CA:918:A:H4'	2.36	0.60
1:CA:984:C:O2'	1:CA:985:C:H5'	2.01	0.60
10:CK:16:SER:HA	10:CK:78:ILE:HA	1.82	0.60
21:CN:5:MET:HG2	21:CN:8:ARG:NE	2.16	0.60
13:CP:67:ILE:HG13	13:CP:71:VAL:HG13	1.84	0.60
23:DB:1360:G:H2'	23:DB:1361:G:H5'	1.84	0.60
23:DB:1788:C:O2'	23:DB:1789:A:H5'	2.01	0.60
23:DB:2859:G:H2'	23:DB:2860:A:C8	2.37	0.60
23:DB:417:C:H2'	23:DB:418:C:C6	2.37	0.60
23:DB:560:C:H2'	23:DB:561:G:O4'	2.01	0.60
23:DB:580:U:H2'	23:DB:581:C:H6	1.67	0.60
23:DB:654:A:H2'	23:DB:654:A:N3	2.16	0.60
23:DB:923:G:H1'	52:DW:23:LYS:HZ2	1.67	0.60
26:DD:149:ASN:O	26:DD:152:PRO:HD2	2.01	0.60
23:DB:1007:C:O3'	41:DJ:110:PRO:HB3	2.02	0.60
43:DO:11:ALA:HB2	43:DO:96:GLY:N	2.17	0.60
1:AA:1513:A:H2'	1:AA:1514:G:C8	2.36	0.60
21:AN:9:GLU:O	21:AN:13:VAL:HG23	2.01	0.60
16:AS:44:ILE:HA	16:AS:61:VAL:HB	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AK:124:LYS:O	19:AU:33:ARG:NE	2.34	0.60
23:BB:2348:U:H5'	33:B1:20:TYR:OH	2.02	0.60
23:BB:528:A:C2	23:BB:2042:A:H2'	2.37	0.60
23:BB:2229:U:H2'	23:BB:2230:G:C8	2.37	0.60
26:BD:168:GLU:O	26:BD:170:VAL:HG22	2.02	0.60
29:BE:145:ASP:N	29:BE:166:LYS:HB3	2.17	0.60
41:BJ:56:VAL:HG12	41:BJ:57:LEU:H	1.66	0.60
27:BK:54:LYS:HD2	27:BK:54:LYS:N	2.16	0.60
28:BP:61:ARG:HH21	28:BP:61:ARG:HB3	1.67	0.60
52:BW:23:LYS:C	52:BW:66:VAL:HB	2.22	0.60
1:CA:1013:G:N2	1:CA:1015:G:H3'	2.17	0.60
2:CC:109:GLU:HB3	2:CC:139:ASN:HB3	1.83	0.60
2:CC:26:LYS:HE2	2:CC:27:GLU:N	2.16	0.60
8:CI:18:VAL:HG21	8:CI:81:GLY:C	2.22	0.60
33:D1:7:LYS:HD3	34:D3:33:THR:HG21	1.82	0.60
23:DB:1551:A:C3'	23:DB:1552:A:H5''	2.31	0.60
23:DB:181:A:H2'	23:DB:182:A:H8	1.65	0.60
23:DB:1897:G:O2'	23:DB:1898:U:H5'	2.01	0.60
23:DB:1914:C:H2'	23:DB:1915:U:C6	2.37	0.60
23:DB:2591:C:H2'	23:DB:2592:G:H8	1.66	0.60
25:DC:155:ARG:CB	25:DC:155:ARG:HH11	2.15	0.60
47:DF:24:VAL:O	47:DF:27:VAL:HG22	2.02	0.60
48:DG:83:THR:HA	48:DG:84:LYS:NZ	2.17	0.60
38:DM:35:ALA:HB3	38:DM:99:GLY:N	2.16	0.60
45:DS:3:THR:HB	45:DS:62:ASP:HB2	1.84	0.60
51:DZ:40:VAL:HG22	51:DZ:45:ARG:O	2.00	0.60
1:AA:939:G:O2'	1:AA:1375:A:H1'	2.02	0.60
1:AA:553:A:H2'	1:AA:554:A:C8	2.37	0.60
1:AA:806:C:H2'	1:AA:807:A:H8	1.67	0.60
5:AF:86:ARG:NH1	15:AR:63:TYR:HB3	2.17	0.60
23:BB:1440:U:H2'	23:BB:1441:G:C8	2.36	0.60
23:BB:2443:C:H2'	23:BB:2444:G:H8	1.66	0.60
23:BB:870:U:H2'	23:BB:871:U:H5''	1.84	0.60
23:BB:910:A:H62	38:BM:12:MET:HA	1.66	0.60
25:BC:124:LYS:HE2	25:BC:127:ASN:HD21	1.65	0.60
26:BD:34:VAL:HG12	26:BD:94:GLN:H	1.64	0.60
27:BK:10:VAL:HG21	27:BK:16:ALA:HA	1.84	0.60
38:BM:41:LEU:HA	38:BM:45:GLN:OE1	2.02	0.60
38:BM:66:ARG:NE	38:BM:101:VAL:HG11	2.17	0.60
49:BR:48:LYS:HD2	49:BR:48:LYS:O	2.02	0.60
45:BS:88:ARG:HH21	45:BS:88:ARG:HG3	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:4:GLU:OE2	50:BT:5:GLU:HG2	2.02	0.60
52:BW:32:ALA:C	52:BW:34:SER:H	2.05	0.60
51:BZ:59:ILE:HD13	51:BZ:67:VAL:HG21	1.84	0.60
1:CA:408:A:H3'	1:CA:409:U:H6	1.67	0.60
9:CJ:53:ILE:CG2	9:CJ:61:ALA:HB1	2.31	0.60
10:CK:28:ASN:HD21	10:CK:46:ALA:HB3	1.65	0.60
10:CK:92:ARG:NH2	19:CU:19:LYS:HG2	2.17	0.60
12:CM:106:ARG:HB3	12:CM:111:PRO:HA	1.84	0.60
33:D1:3:GLY:O	33:D1:4:ILE:HG12	2.02	0.60
36:D2:20:ALA:C	36:D2:22:MET:H	2.04	0.60
22:DA:6:G:H2'	22:DA:7:G:H8	1.66	0.60
23:DB:2262:U:H2'	23:DB:2263:C:H6	1.67	0.60
23:DB:2286:G:H4'	23:DB:2287:A:O4'	2.01	0.60
23:DB:551:G:H2'	23:DB:552:U:C6	2.37	0.60
23:DB:919:U:H2'	23:DB:920:A:H8	1.67	0.60
26:DD:37:VAL:HG23	26:DD:91:THR:HA	1.84	0.60
29:DE:145:ASP:HB3	29:DE:184:ASP:H	1.66	0.60
29:DE:1:MET:HB3	29:DE:14:VAL:HG23	1.83	0.60
23:DB:659:G:H21	29:DE:30:GLN:NE2	2.00	0.60
44:DQ:91:ARG:HB2	44:DQ:94:LEU:HD23	1.83	0.60
49:DR:49:ILE:HD12	49:DR:49:ILE:O	2.02	0.60
46:DU:85:ARG:NH1	46:DU:86:PHE:H	2.00	0.60
52:DW:23:LYS:C	52:DW:66:VAL:HB	2.22	0.60
1:AA:944:G:H21	1:AA:1339:A:H62	1.50	0.59
1:AA:373:A:H2'	1:AA:374:A:H8	1.67	0.59
1:AA:468:A:H3'	1:AA:469:C:C6	2.37	0.59
18:AB:79:VAL:HG12	18:AB:90:PHE:HB2	1.83	0.59
3:AD:58:GLN:O	3:AD:62:ARG:HG2	2.02	0.59
9:AJ:10:LEU:O	9:AJ:71:LEU:HB2	2.02	0.59
36:B2:19:ARG:NH2	36:B2:19:ARG:HB3	2.17	0.59
23:BB:1353:A:H2'	23:BB:1354:A:C8	2.36	0.59
23:BB:2229:U:H2'	23:BB:2230:G:H8	1.67	0.59
23:BB:2365:G:O2'	52:BW:59:PHE:HE1	1.85	0.59
23:BB:642:U:O2	23:BB:644:A:H3'	2.02	0.59
26:BD:170:VAL:O	26:BD:170:VAL:HG23	2.01	0.59
26:BD:33:ARG:NH1	26:BD:51:THR:HG22	2.16	0.59
23:BB:321:U:H1'	29:BE:162:ARG:NH1	2.14	0.59
29:BE:52:VAL:HG12	29:BE:53:THR:N	2.14	0.59
29:BE:60:TRP:CH2	29:BE:69:ARG:HA	2.37	0.59
27:BK:102:PRO:HA	27:BK:120:PRO:HB3	1.84	0.59
27:BK:86:LEU:H	27:BK:86:LEU:HD23	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BL:2:ARG:HG2	37:BL:5:THR:HG21	1.82	0.59
44:BQ:42:GLY:HA3	49:BR:75:VAL:HG21	1.82	0.59
46:BU:64:ILE:HD11	46:BU:68:ASN:HD22	1.67	0.59
52:BW:40:ARG:O	52:BW:44:PHE:HA	2.01	0.59
1:CA:844:G:N2	1:CA:845:A:H62	2.00	0.59
3:CD:29:THR:HG22	3:CD:30:LYS:N	2.17	0.59
5:CF:53:LYS:HB2	5:CF:54:LEU:HD22	1.82	0.59
9:CJ:36:VAL:HA	9:CJ:77:VAL:HG23	1.84	0.59
19:CU:35:GLU:HB2	19:CU:37:TYR:CE2	2.36	0.59
32:D4:1:MET:HE1	32:D4:36:ARG:HB2	1.84	0.59
23:DB:2810:A:H2'	23:DB:2811:G:O4'	2.02	0.59
23:DB:290:U:O2'	23:DB:291:G:H5'	2.02	0.59
23:DB:581:C:H2'	23:DB:582:A:C8	2.37	0.59
26:DD:168:GLU:HG3	26:DD:170:VAL:HG13	1.83	0.59
26:DD:39:ASP:HB3	26:DD:42:ASN:HB3	1.83	0.59
29:DE:188:MET:HG3	29:DE:192:ALA:HB3	1.83	0.59
27:DK:10:VAL:HG21	27:DK:16:ALA:HA	1.83	0.59
27:DK:25:LEU:HB2	27:DK:38:ILE:HG12	1.84	0.59
37:DL:93:ASN:O	37:DL:95:LEU:N	2.35	0.59
46:DU:32:LYS:HA	46:DU:65:GLN:HA	1.84	0.59
23:DB:855:G:N3	52:DW:23:LYS:HE3	2.17	0.59
23:DB:928:A:O2'	30:DY:37:ARG:HD3	2.02	0.59
51:DZ:21:ALA:HB3	51:DZ:23:ASN:ND2	2.17	0.59
1:AA:713:G:H2'	1:AA:714:G:H8	1.66	0.59
1:AA:93:U:O5'	1:AA:93:U:H6	1.85	0.59
1:AA:971:G:H3'	1:AA:971:G:OP1	2.02	0.59
18:AB:209:VAL:O	18:AB:213:LEU:HB2	2.02	0.59
8:AI:17:ARG:HH21	8:AI:65:THR:HG21	1.66	0.59
8:AI:27:ILE:HB	8:AI:34:LEU:HD22	1.84	0.59
10:AK:70:ALA:HA	10:AK:73:VAL:HG22	1.83	0.59
19:AU:35:GLU:HB2	19:AU:37:TYR:CE2	2.36	0.59
36:B2:31:LEU:HD22	36:B2:42:LEU:HD12	1.83	0.59
23:BB:1722:A:H2'	23:BB:1723:G:H8	1.68	0.59
23:BB:208:C:H2'	23:BB:209:C:H6	1.67	0.59
23:BB:2186:G:H2'	23:BB:2187:U:O4'	2.02	0.59
23:BB:2849:U:H4'	23:BB:2850:A:H5'	1.83	0.59
23:BB:506:G:H5'	23:BB:509:C:H1'	1.84	0.59
23:BB:802:A:H2'	23:BB:803:U:C6	2.37	0.59
25:BC:202:ARG:O	25:BC:203:VAL:HB	2.01	0.59
26:BD:104:VAL:HG12	26:BD:106:LYS:HE2	1.84	0.59
26:BD:174:SER:O	26:BD:175:LEU:HB2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2787:C:H5'	26:BD:66:GLY:HA3	1.83	0.59
47:BF:120:SER:HG	47:BF:127:TYR:HD2	1.50	0.59
49:BR:24:LYS:HA	49:BR:94:THR:OG1	2.01	0.59
50:BT:69:ARG:HG2	50:BT:74:ILE:N	2.17	0.59
46:BU:54:PRO:HG2	46:BU:55:GLY:H	1.65	0.59
1:CA:1141:C:H2'	1:CA:1142:G:C8	2.37	0.59
1:CA:1170:A:H2'	1:CA:1171:A:O4'	2.01	0.59
1:CA:1412:C:H2'	1:CA:1413:A:H8	1.68	0.59
18:CB:86:CYS:SG	18:CB:88:GLN:HG3	2.42	0.59
6:CG:70:PRO:HB3	6:CG:102:TRP:CH2	2.38	0.59
7:CH:125:ILE:HG22	7:CH:126:CYS:SG	2.42	0.59
7:CH:73:SER:O	7:CH:128:VAL:HA	2.02	0.59
1:CA:255:G:H4'	14:CQ:18:LYS:HD2	1.84	0.59
17:CT:43:LYS:HZ2	17:CT:43:LYS:N	2.00	0.59
22:DA:32:U:H1'	22:DA:52:A:N7	2.17	0.59
23:DB:1475:G:H4'	23:DB:1476:U:O5'	2.02	0.59
23:DB:160:A:H2'	23:DB:161:A:C8	2.36	0.59
23:DB:2292:U:H2'	23:DB:2293:G:H8	1.67	0.59
23:DB:797:G:OP2	29:DE:57:LYS:HB2	2.02	0.59
47:DF:134:GLN:NE2	47:DF:136:ILE:HD13	2.17	0.59
48:DG:25:ILE:O	48:DG:32:LEU:HA	2.02	0.59
27:DK:59:LYS:HD2	27:DK:89:ASN:HD22	1.68	0.59
38:DM:12:MET:HB2	38:DM:72:PRO:HG2	1.84	0.59
23:DB:2334:U:H5'	43:DO:12:THR:HB	1.83	0.59
46:DU:47:PRO:HD3	46:DU:55:GLY:HA3	1.83	0.59
39:DX:23:ARG:H	39:DX:26:PHE:HB3	1.67	0.59
2:AC:127:VAL:HG23	2:AC:128:MET:N	2.17	0.59
4:AE:148:SER:HB2	4:AE:149:PRO:HD2	1.84	0.59
1:AA:1180:A:P	8:AI:98:ARG:HH22	2.26	0.59
11:AL:34:THR:HB	11:AL:53:ARG:HB3	1.82	0.59
23:BB:1021:A:H2'	23:BB:1023:U:H5''	1.85	0.59
23:BB:129:C:H2'	23:BB:130:C:H6	1.65	0.59
23:BB:1447:C:H2'	23:BB:1448:G:H8	1.67	0.59
23:BB:1580:A:H2'	23:BB:1581:G:O4'	2.01	0.59
23:BB:20:C:H2'	23:BB:21:A:C8	2.36	0.59
23:BB:2243:U:O2'	23:BB:2244:U:H5'	2.00	0.59
29:BE:188:MET:HG3	29:BE:192:ALA:HB3	1.83	0.59
47:BF:139:GLU:HG2	47:BF:140:ILE:H	1.67	0.59
48:BG:15:ASP:HB3	48:BG:26:LYS:H	1.66	0.59
48:BG:24:THR:HA	48:BG:33:THR:O	2.02	0.59
40:BH:2:GLN:O	40:BH:3:VAL:HG22	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:73:ASN:ND2	40:BH:74:ALA:H	2.01	0.59
45:BS:46:LEU:O	45:BS:50:VAL:HG23	2.02	0.59
35:BV:72:VAL:CG1	35:BV:93:ARG:HA	2.33	0.59
1:CA:1504:G:H4'	1:CA:1505:G:O4'	2.01	0.59
1:CA:60:A:H4'	1:CA:61:G:OP1	2.02	0.59
18:CB:65:LYS:HB2	18:CB:158:ASP:OD1	2.01	0.59
2:CC:122:GLN:HB3	2:CC:127:VAL:HG21	1.83	0.59
2:CC:63:ILE:O	2:CC:98:ALA:HB1	2.02	0.59
4:CE:148:SER:HB2	4:CE:149:PRO:HD2	1.82	0.59
6:CG:43:TYR:O	6:CG:47:GLU:HG2	2.02	0.59
1:CA:135:C:N3	13:CP:1:MET:HB2	2.18	0.59
22:DA:75:G:H2'	22:DA:76:G:H8	1.66	0.59
23:DB:1082:U:C2	23:DB:1086:A:C6	2.91	0.59
23:DB:1631:G:N2	23:DB:1633:G:H3'	2.17	0.59
23:DB:1635:A:C2'	23:DB:1636:U:H5'	2.32	0.59
23:DB:2784:U:H2'	23:DB:2785:C:H6	1.66	0.59
23:DB:834:G:H2'	23:DB:835:C:O4'	2.01	0.59
27:DK:17:ARG:HB2	27:DK:45:GLU:HB3	1.84	0.59
1:AA:1043:G:H2'	1:AA:1044:A:H8	1.67	0.59
1:AA:517:G:H22	1:AA:533:A:P	2.26	0.59
1:AA:810:C:O2'	1:AA:811:C:H5'	2.02	0.59
18:AB:19:THR:HG23	18:AB:20:ARG:H	1.66	0.59
5:AF:86:ARG:CZ	15:AR:63:TYR:HB3	2.32	0.59
17:AT:48:LYS:HG3	17:AT:49:ALA:N	2.18	0.59
23:BB:1166:G:H2'	23:BB:1167:C:C6	2.37	0.59
23:BB:2085:U:O2'	23:BB:2086:U:H5'	2.01	0.59
23:BB:2630:G:H2'	23:BB:2631:G:H8	1.67	0.59
25:BC:94:LEU:HB2	25:BC:100:ARG:HB3	1.85	0.59
23:BB:558:U:H5''	41:BJ:111:LYS:HD3	1.84	0.59
42:BN:94:TYR:O	42:BN:115:LEU:HA	2.02	0.59
49:BR:39:LEU:CB	49:BR:49:ILE:HD11	2.32	0.59
45:BS:10:ALA:HB3	45:BS:101:SER:OG	2.01	0.59
46:BU:27:VAL:HG23	46:BU:33:VAL:HG12	1.83	0.59
1:CA:1306:A:H2'	1:CA:1307:U:O4'	2.02	0.59
8:CI:42:THR:HA	8:CI:45:MET:SD	2.42	0.59
10:CK:16:SER:HA	10:CK:77:GLY:O	2.03	0.59
32:D4:24:ARG:HG2	32:D4:36:ARG:HG3	1.83	0.59
23:DB:1100:C:H2'	23:DB:1101:U:H6	1.67	0.59
23:DB:1440:U:H2'	23:DB:1441:G:C8	2.37	0.59
23:DB:2041:U:H2'	23:DB:2042:A:C8	2.37	0.59
23:DB:528:A:C2	23:DB:2042:A:H2'	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:43:ASN:HB2	25:DC:49:THR:HG23	1.84	0.59
26:DD:170:VAL:O	26:DD:170:VAL:HG23	2.01	0.59
47:DF:36:ASN:HA	47:DF:87:LYS:HA	1.84	0.59
40:DH:2:GLN:O	40:DH:3:VAL:HG22	2.02	0.59
28:DP:31:VAL:O	28:DP:32:VAL:HB	2.03	0.59
28:DP:31:VAL:HG12	28:DP:38:ARG:O	2.03	0.59
52:DW:44:PHE:O	52:DW:78:PHE:HA	2.02	0.59
52:DW:77:LYS:O	52:DW:78:PHE:HB2	2.00	0.59
1:AA:1175:G:O2'	1:AA:1176:A:H5'	2.02	0.59
1:AA:628:G:H2'	1:AA:629:A:H8	1.67	0.59
1:AA:74:A:H2'	1:AA:75:G:O4'	2.01	0.59
1:AA:812:G:H2'	1:AA:812:G:N3	2.15	0.59
2:AC:6:PRO:HG2	2:AC:183:TYR:CD2	2.37	0.59
7:AH:94:VAL:HG12	7:AH:99:GLY:HA3	1.84	0.59
12:AM:53:ASP:HA	12:AM:56:ARG:HH21	1.68	0.59
10:AK:126:ARG:HB2	19:AU:33:ARG:HD2	1.82	0.59
19:AU:48:LYS:HA	19:AU:51:ALA:HB3	1.84	0.59
22:BA:41:G:O6	47:BF:68:LYS:HD3	2.03	0.59
23:BB:2087:G:H2'	23:BB:2088:A:H8	1.67	0.59
23:BB:275:C:H2'	23:BB:276:U:H5'	1.83	0.59
23:BB:417:C:H2'	23:BB:418:C:C6	2.37	0.59
41:BJ:30:THR:N	41:BJ:108:MET:HE3	2.18	0.59
45:BS:81:SER:HA	45:BS:99:ARG:HA	1.84	0.59
46:BU:34:ILE:HG12	46:BU:63:ALA:HB2	1.83	0.59
46:BU:47:PRO:HD3	46:BU:55:GLY:HA3	1.85	0.59
52:BW:77:LYS:O	52:BW:78:PHE:HB2	2.01	0.59
30:BY:26:LEU:HD12	30:BY:28:LEU:HD11	1.85	0.59
1:CA:190:A:O5'	1:CA:190:A:H8	1.85	0.59
1:CA:98:A:H2'	1:CA:99:C:H6	1.67	0.59
3:CD:187:ARG:NH1	3:CD:191:SER:HA	2.18	0.59
7:CH:55:LYS:HE3	7:CH:55:LYS:HA	1.84	0.59
5:CF:86:ARG:NH1	15:CR:63:TYR:HB3	2.17	0.59
23:DB:1021:A:H61	23:DB:1142:A:N6	2.00	0.59
23:DB:1549:A:H2'	23:DB:1550:C:C6	2.38	0.59
23:DB:2384:U:H5''	23:DB:2386:A:OP1	2.03	0.59
25:DC:89:ASN:O	25:DC:105:ALA:HB3	2.02	0.59
40:DH:2:GLN:HB3	40:DH:18:GLN:HG3	1.83	0.59
41:DJ:29:ALA:HA	41:DJ:32:LEU:HB2	1.83	0.59
37:DL:109:LYS:HB2	37:DL:111:ILE:CD1	2.32	0.59
44:DQ:77:LYS:HA	44:DQ:80:ASN:HB3	1.85	0.59
44:DQ:111:LYS:HB2	49:DR:48:LYS:NZ	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:43:ILE:O	50:DT:47:VAL:HG23	2.03	0.59
1:AA:188:C:H2'	1:AA:189:A:O4'	2.03	0.59
1:AA:638:U:H2'	1:AA:639:G:O4'	2.02	0.59
3:AD:59:LYS:O	3:AD:63:ILE:HG13	2.02	0.59
4:AE:89:THR:HG21	4:AE:134:ASN:HD21	1.66	0.59
5:AF:73:GLU:O	5:AF:77:THR:HG23	2.03	0.59
6:AG:99:ALA:O	6:AG:103:ILE:HG13	2.03	0.59
12:AM:19:THR:HA	12:AM:24:VAL:HG23	1.83	0.59
23:BB:1046:A:C4'	23:BB:1047:G:H5''	2.31	0.59
23:BB:1552:A:H2'	23:BB:1553:A:H5'	1.85	0.59
23:BB:438:G:H2'	23:BB:439:A:H8	1.67	0.59
23:BB:547:A:H2'	23:BB:547:A:N3	2.18	0.59
1:CA:1266:G:N2	1:CA:1268:G:H3'	2.18	0.59
1:CA:937:A:H1'	1:CA:1379:G:N2	2.18	0.59
1:CA:551:U:O2'	11:CL:82:ARG:HD2	2.02	0.59
1:CA:557:G:H2'	1:CA:558:G:O4'	2.02	0.59
1:CA:638:U:H2'	1:CA:639:G:O4'	2.02	0.59
1:CA:953:G:H2'	1:CA:954:G:O4'	2.02	0.59
18:CB:169:HIS:HA	18:CB:172:ILE:HD12	1.83	0.59
8:CI:54:VAL:HG11	8:CI:86:LEU:HD23	1.85	0.59
21:CN:25:GLU:O	21:CN:29:ILE:HG13	2.03	0.59
20:CO:32:LEU:HD13	20:CO:35:GLN:HE22	1.66	0.59
32:D4:16:ILE:HG12	32:D4:25:VAL:HG22	1.84	0.59
23:DB:136:G:H2'	23:DB:137:U:C6	2.37	0.59
23:DB:2019:A:H2	23:DB:2035:G:H22	1.51	0.59
23:DB:918:A:H2'	23:DB:919:U:H5'	1.83	0.59
25:DC:94:LEU:HA	25:DC:100:ARG:HA	1.83	0.59
29:DE:58:LYS:HE2	29:DE:60:TRP:CD1	2.27	0.59
24:DI:112:LYS:O	24:DI:116:MET:HG3	2.02	0.59
27:DK:87:LEU:HB2	27:DK:93:GLN:O	2.01	0.59
38:DM:41:LEU:HA	38:DM:45:GLN:OE1	2.02	0.59
38:DM:61:GLY:HA2	38:DM:107:GLY:HA3	1.84	0.59
42:DN:103:ARG:HB2	42:DN:110:MET:HE2	1.83	0.59
43:DO:58:ILE:O	43:DO:62:LEU:HG	2.01	0.59
44:DQ:80:ASN:O	44:DQ:83:LYS:HB3	2.01	0.59
52:DW:59:PHE:O	52:DW:60:ALA:HB3	2.03	0.59
1:AA:1458:G:H2'	1:AA:1459:G:H8	1.67	0.59
1:AA:71:A:H61	1:AA:99:C:C1'	2.08	0.59
3:AD:151:GLN:HB3	3:AD:153:ARG:HG2	1.84	0.59
5:AF:79:ARG:NH2	5:AF:87:SER:HB3	2.17	0.59
10:AK:16:SER:HA	10:AK:77:GLY:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AS:28:LYS:HD2	16:AS:28:LYS:H	1.67	0.59
33:B1:3:GLY:C	33:B1:5:ARG:H	2.06	0.59
22:BA:55:U:H2'	22:BA:56:G:C8	2.37	0.59
23:BB:1360:G:H2'	23:BB:1361:G:H5'	1.84	0.59
23:BB:1387:A:C5'	23:BB:1469:A:H1'	2.33	0.59
23:BB:1536:C:H4'	23:BB:1537:G:C4	2.37	0.59
23:BB:2462:C:H2'	23:BB:2463:C:C6	2.37	0.59
23:BB:2569:G:O2'	23:BB:2570:G:H5'	2.02	0.59
23:BB:2671:G:H2'	23:BB:2672:U:C6	2.38	0.59
23:BB:971:G:H2'	23:BB:972:A:O4'	2.03	0.59
25:BC:103:ILE:HG22	25:BC:105:ALA:H	1.67	0.59
26:BD:37:VAL:HG23	26:BD:91:THR:HA	1.83	0.59
47:BF:37:MET:HG3	47:BF:150:GLY:O	2.02	0.59
47:BF:94:ARG:HA	47:BF:97:GLU:HB3	1.83	0.59
40:BH:122:LEU:HD12	40:BH:122:LEU:H	1.67	0.59
28:BP:26:GLU:HG3	28:BP:43:GLU:HB2	1.84	0.59
28:BP:13:LYS:HD2	28:BP:76:HIS:HA	1.84	0.59
44:BQ:88:GLU:HG2	49:BR:49:ILE:O	2.03	0.59
52:BW:17:ALA:HA	52:BW:35:ILE:CG2	2.26	0.59
52:BW:59:PHE:O	52:BW:60:ALA:HB3	2.02	0.59
5:CF:91:ARG:H	5:CF:93:LYS:HZ1	1.49	0.59
1:CA:707:U:H4'	10:CK:21:HIS:CD2	2.38	0.59
21:CN:11:LYS:O	21:CN:15:LEU:HG	2.02	0.59
23:DB:21:A:H2'	23:DB:22:C:C6	2.37	0.59
23:DB:2529:G:H4'	48:DG:174:LYS:HE2	1.85	0.59
23:DB:654:A:H2'	23:DB:655:A:H5''	1.84	0.59
23:DB:654:A:N3	23:DB:655:A:H5''	2.17	0.59
23:DB:813:U:H2'	23:DB:814:C:H6	1.68	0.59
25:DC:124:LYS:HE2	25:DC:127:ASN:ND2	2.17	0.59
25:DC:121:ALA:HB3	25:DC:129:LEU:HD21	1.84	0.59
26:DD:174:SER:O	26:DD:175:LEU:HB2	2.02	0.59
26:DD:55:LYS:HZ2	26:DD:55:LYS:HB3	1.68	0.59
47:DF:2:LYS:H	47:DF:2:LYS:CE	2.15	0.59
47:DF:48:LEU:HD23	47:DF:48:LEU:H	1.67	0.59
47:DF:35:LEU:HD22	47:DF:56:LEU:HD11	1.85	0.59
48:DG:104:LEU:HD22	48:DG:106:LEU:HD21	1.84	0.59
40:DH:72:ILE:HG23	40:DH:108:VAL:HG21	1.85	0.59
27:DK:54:LYS:HD2	27:DK:54:LYS:N	2.14	0.59
43:DO:7:ARG:HA	43:DO:10:ARG:NE	2.17	0.59
45:DS:82:MET:HE1	45:DS:84:ARG:HH22	1.67	0.59
50:DT:32:LEU:N	50:DT:83:ALA:HB3	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:34:VAL:HG21	50:DT:43:ILE:HD12	1.84	0.59
1:AA:126:G:H4'	1:AA:634:C:H1'	1.85	0.59
1:AA:1409:C:N3	1:AA:1492:A:H2	2.00	0.59
1:AA:1512:U:H2'	1:AA:1513:A:H8	1.68	0.59
2:AC:111:ASP:OD2	2:AC:114:LEU:HG	2.03	0.59
2:AC:126:ARG:HH11	2:AC:126:ARG:CA	2.15	0.59
4:AE:37:VAL:HA	4:AE:47:PHE:HA	1.85	0.59
6:AG:94:ARG:O	6:AG:98:LEU:HG	2.02	0.59
12:AM:89:ARG:HD2	12:AM:95:PRO:O	2.01	0.59
33:B1:10:LEU:O	33:B1:19:PHE:HB2	2.03	0.59
22:BA:95:U:H2'	22:BA:96:G:C8	2.37	0.59
23:BB:2514:U:H5''	41:BJ:81:ILE:HD12	1.84	0.59
23:BB:2693:G:H2'	23:BB:2694:G:H8	1.68	0.59
47:BF:97:GLU:O	47:BF:100:GLU:HB2	2.01	0.59
40:BH:140:ALA:C	40:BH:141:LYS:HD3	2.23	0.59
41:BJ:13:ARG:O	41:BJ:52:ASP:HA	2.03	0.59
37:BL:143:GLU:HG2	37:BL:144:GLU:N	2.14	0.59
50:BT:32:LEU:HG	50:BT:83:ALA:HB2	1.85	0.59
1:CA:1382:C:H4'	6:CG:78:ARG:NH2	2.17	0.59
1:CA:235:C:H2'	1:CA:236:A:H8	1.67	0.59
1:CA:411:A:H2'	1:CA:412:A:N3	2.17	0.59
1:CA:469:C:H2'	1:CA:470:C:C6	2.38	0.59
1:CA:513:C:H2'	1:CA:514:C:H6	1.67	0.59
1:CA:711:G:O2'	1:CA:712:A:H5'	2.02	0.59
1:CA:900:A:H2'	1:CA:901:A:H8	1.64	0.59
1:CA:951:G:H1'	1:CA:970:C:O2'	2.02	0.59
18:CB:135:MET:O	18:CB:138:ARG:HB3	2.03	0.59
4:CE:71:ILE:HG12	4:CE:72:ASN:N	2.16	0.59
8:CI:35:GLU:O	8:CI:39:GLY:HA3	2.01	0.59
33:D1:3:GLY:C	33:D1:5:ARG:H	2.06	0.59
22:DA:32:U:C4'	22:DA:52:A:H62	2.15	0.59
23:DB:1177:G:H2'	23:DB:1178:C:H6	1.67	0.59
23:DB:1387:A:C5'	23:DB:1469:A:H1'	2.31	0.59
23:DB:1443:U:H2'	23:DB:1444:G:H8	1.68	0.59
23:DB:2261:C:H1'	23:DB:2388:A:N3	2.16	0.59
23:DB:2557:G:H2'	23:DB:2558:C:C6	2.37	0.59
23:DB:63:A:H2'	23:DB:64:A:N7	2.17	0.59
26:DD:33:ARG:HH11	26:DD:51:THR:CG2	2.15	0.59
28:DP:19:PHE:HB2	28:DP:50:ARG:HH12	1.67	0.59
44:DQ:103:VAL:O	44:DQ:106:THR:HB	2.03	0.59
1:AA:1075:U:H2'	1:AA:1076:U:C6	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:202:G:H2'	1:AA:203:G:H8	1.67	0.59
1:AA:328:C:H4'	1:AA:329:A:C5'	2.33	0.59
1:AA:920:U:H2'	1:AA:921:U:H6	1.68	0.59
4:AE:106:ALA:HB1	4:AE:110:MET:HB3	1.85	0.59
7:AH:31:LEU:HG	7:AH:35:ILE:HD11	1.85	0.59
23:BB:1175:A:H3'	23:BB:1176:U:H4'	1.85	0.59
23:BB:1874:C:H2'	23:BB:1875:G:O4'	2.03	0.59
23:BB:532:A:N1	23:BB:2020:A:H1'	2.17	0.59
23:BB:2728:U:H2'	23:BB:2729:G:H8	1.67	0.59
26:BD:33:ARG:NE	26:BD:51:THR:HB	2.17	0.59
40:BH:99:ILE:HG21	40:BH:130:VAL:CB	2.32	0.59
27:BK:43:ILE:HD12	27:BK:56:ASP:HB2	1.85	0.59
43:BO:115:LEU:C	43:BO:116:GLN:HE21	2.05	0.59
46:BU:42:LYS:HD2	46:BU:59:GLU:HB2	1.83	0.59
1:CA:1232:U:H5''	8:CI:125:GLN:O	2.02	0.59
1:CA:673:A:H2'	1:CA:674:G:C8	2.37	0.59
18:CB:134:LEU:HD12	18:CB:135:MET:N	2.18	0.59
18:CB:95:TRP:CZ3	18:CB:99:MET:HG3	2.38	0.59
3:CD:58:GLN:O	3:CD:62:ARG:HG2	2.02	0.59
21:CN:72:PHE:O	21:CN:73:LEU:HD23	2.03	0.59
14:CQ:75:VAL:HG23	14:CQ:76:ARG:H	1.68	0.59
19:CU:48:LYS:HA	19:CU:51:ALA:HB3	1.84	0.59
23:DB:1354:A:H2'	23:DB:1355:G:O4'	2.03	0.59
23:DB:2400:G:O2'	23:DB:2401:U:H5'	2.02	0.59
23:DB:480:A:H3'	23:DB:481:G:H5''	1.85	0.59
23:DB:642:U:O2	23:DB:644:A:H3'	2.03	0.59
25:DC:20:ASN:ND2	25:DC:23:LEU:HD13	2.18	0.59
25:DC:94:LEU:HB2	25:DC:100:ARG:HB3	1.85	0.59
40:DH:58:LEU:O	40:DH:61:VAL:HG12	2.02	0.59
28:DP:61:ARG:HH21	28:DP:61:ARG:HB3	1.68	0.59
50:DT:9:LYS:NZ	50:DT:9:LYS:HB3	2.18	0.59
35:DV:63:ILE:N	35:DV:63:ILE:HD12	2.18	0.59
52:DW:40:ARG:O	52:DW:44:PHE:HA	2.02	0.59
30:DY:15:ARG:O	30:DY:20:LYS:HE3	2.03	0.59
18:AB:119:GLN:O	18:AB:125:PHE:HB3	2.01	0.59
18:AB:26:MET:HG3	18:AB:188:THR:O	2.03	0.59
4:AE:71:ILE:HG12	4:AE:72:ASN:N	2.15	0.59
4:AE:95:MET:HA	4:AE:124:ALA:HB2	1.85	0.59
6:AG:71:THR:O	6:AG:89:GLU:HA	2.03	0.59
7:AH:45:ILE:HA	7:AH:63:LYS:HB2	1.84	0.59
8:AI:61:ASP:C	8:AI:62:LEU:HD13	2.23	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AK:108:ASN:ND2	19:AU:6:ARG:HD2	2.17	0.59
11:AL:98:ARG:HH21	11:AL:104:SER:C	2.06	0.59
1:AA:1458:G:H4'	17:AT:22:SER:HB2	1.84	0.59
22:BA:6:G:H2'	22:BA:7:G:C8	2.37	0.59
23:BB:1106:G:H2'	23:BB:1107:G:H8	1.67	0.59
23:BB:131:A:O2'	23:BB:132:G:H5'	2.01	0.59
23:BB:1549:A:H2'	23:BB:1550:C:C6	2.38	0.59
23:BB:2567:G:H2'	23:BB:2568:U:C6	2.37	0.59
23:BB:2784:U:H2'	23:BB:2785:C:H6	1.67	0.59
24:BI:96:LYS:N	24:BI:96:LYS:HD2	2.18	0.59
44:BQ:103:VAL:O	44:BQ:106:THR:HB	2.03	0.59
31:B0:21:LEU:HD12	45:BS:19:LEU:O	2.03	0.59
45:BS:3:THR:HB	45:BS:62:ASP:HB2	1.84	0.59
1:CA:430:A:OP1	3:CD:8:LEU:HB2	2.03	0.59
1:CA:556:C:O2'	1:CA:557:G:H5'	2.03	0.59
18:CB:69:VAL:HB	18:CB:162:VAL:CG2	2.33	0.59
8:CI:94:ARG:HE	8:CI:97:LEU:HD12	1.68	0.59
14:CQ:16:MET:CB	14:CQ:19:SER:HB2	2.33	0.59
16:CS:10:ILE:CG2	16:CS:37:SER:HB3	2.32	0.59
17:CT:79:THR:O	17:CT:82:ILE:HG13	2.03	0.59
31:D0:53:VAL:HB	42:DN:118:ARG:HH22	1.67	0.59
22:DA:6:G:H2'	22:DA:7:G:C8	2.38	0.59
23:DB:1537:G:H3'	23:DB:1537:G:N3	2.18	0.59
23:DB:1580:A:H2'	23:DB:1581:G:O4'	2.03	0.59
23:DB:710:U:H2'	23:DB:711:G:H8	1.67	0.59
27:DK:3:GLN:HG2	27:DK:4:GLU:N	2.18	0.59
37:DL:68:SER:HB2	37:DL:71:ALA:HB3	1.84	0.59
44:DQ:108:LEU:HA	49:DR:48:LYS:HD3	1.84	0.59
35:DV:62:THR:HG22	35:DV:71:LYS:HG2	1.85	0.59
1:AA:1168:U:H4'	1:AA:1169:A:OP2	2.03	0.58
2:AC:59:PRO:HG2	2:AC:62:SER:OG	2.02	0.58
1:AA:716:A:N3	10:AK:119:GLY:HA2	2.18	0.58
12:AM:87:GLY:HA2	12:AM:90:HIS:HD2	1.67	0.58
16:AS:39:ILE:HB	16:AS:66:VAL:O	2.02	0.58
19:AU:42:THR:O	19:AU:46:ARG:HG3	2.03	0.58
33:B1:31:GLU:CD	33:B1:31:GLU:H	2.05	0.58
23:BB:1443:U:H2'	23:BB:1444:G:H8	1.68	0.58
23:BB:2557:G:H2'	23:BB:2558:C:C6	2.38	0.58
23:BB:519:U:H2'	23:BB:520:G:H8	1.68	0.58
23:BB:856:G:H1'	52:BW:23:LYS:HB3	1.84	0.58
23:BB:912:C:O2'	23:BB:913:U:H5'	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:947:A:H2'	23:BB:948:C:H6	1.67	0.58
25:BC:177:SER:O	25:BC:270:ARG:HG3	2.03	0.58
27:BK:2:ILE:HD12	27:BK:2:ILE:N	2.18	0.58
1:CA:205:A:H2'	1:CA:206:C:H6	1.67	0.58
1:CA:847:G:H2'	1:CA:848:C:C6	2.38	0.58
2:CC:152:VAL:HB	2:CC:165:GLU:HB2	1.85	0.58
3:CD:89:LEU:HD22	3:CD:199:ILE:HD11	1.85	0.58
6:CG:112:ASP:HB2	6:CG:118:ARG:HG2	1.84	0.58
10:CK:28:ASN:CG	10:CK:56:LYS:HE3	2.23	0.58
12:CM:79:LEU:HB2	12:CM:84:CYS:SG	2.42	0.58
23:DB:1166:G:H2'	23:DB:1167:C:C6	2.38	0.58
23:DB:1458:U:H5''	23:DB:1459:G:OP1	2.03	0.58
23:DB:1485:U:H2'	23:DB:1486:U:H6	1.68	0.58
23:DB:532:A:N1	23:DB:2020:A:H1'	2.18	0.58
23:DB:2773:C:H5''	26:DD:169:ARG:HB2	1.84	0.58
25:DC:221:GLY:O	25:DC:223:ALA:N	2.30	0.58
29:DE:21:ARG:HG3	29:DE:22:ASP:N	2.18	0.58
23:DB:471:A:H5''	29:DE:79:ARG:HH12	1.67	0.58
23:DB:1060:U:C5	24:DI:131:THR:HG22	2.38	0.58
43:DO:58:ILE:HG22	43:DO:62:LEU:HD21	1.86	0.58
1:AA:190:A:O5'	1:AA:190:A:H8	1.84	0.58
1:AA:551:U:O2'	11:AL:82:ARG:HD2	2.03	0.58
1:AA:858:G:O6	1:AA:869:G:H3'	2.03	0.58
2:AC:179:ALA:HB1	2:AC:202:PHE:CE1	2.39	0.58
3:AD:187:ARG:NH1	3:AD:191:SER:HA	2.18	0.58
6:AG:108:ARG:HG2	6:AG:115:MET:SD	2.42	0.58
9:AJ:7:ARG:O	9:AJ:100:ILE:HB	2.03	0.58
1:AA:1225:A:OP1	12:AM:100:ARG:HA	2.03	0.58
23:BB:1414:C:H2'	23:BB:1415:U:C6	2.39	0.58
25:BC:89:ASN:O	25:BC:105:ALA:HB3	2.03	0.58
27:BK:25:LEU:HD21	27:BK:40:LYS:N	2.19	0.58
42:BN:33:ILE:HB	42:BN:114:GLU:HB2	1.84	0.58
43:BO:28:VAL:HG11	43:BO:92:PHE:CZ	2.39	0.58
35:BV:61:LEU:O	35:BV:71:LYS:HA	2.03	0.58
1:CA:1343:G:H2'	1:CA:1344:C:C6	2.38	0.58
1:CA:209:U:H5'	1:CA:210:C:C5	2.37	0.58
1:CA:845:A:H2'	1:CA:845:A:N3	2.18	0.58
18:CB:62:ARG:H	18:CB:62:ARG:HD2	1.68	0.58
18:CB:83:ALA:HB3	18:CB:90:PHE:HB3	1.85	0.58
9:CJ:12:ALA:HB1	9:CJ:17:LEU:HD12	1.84	0.58
9:CJ:53:ILE:HG23	9:CJ:61:ALA:HB1	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:40:ILE:O	9:CJ:72:ARG:HA	2.02	0.58
12:CM:104:ASN:H	12:CM:106:ARG:NH2	1.97	0.58
21:CN:81:ILE:HD12	21:CN:82:LYS:H	1.67	0.58
23:DB:2284:A:OP2	33:D1:5:ARG:HG3	2.02	0.58
23:DB:1092:C:H2'	23:DB:1093:G:H5'	1.85	0.58
23:DB:1149:G:H2'	23:DB:1150:C:H6	1.68	0.58
23:DB:1868:C:H2'	23:DB:1869:G:O4'	2.04	0.58
23:DB:20:C:H2'	23:DB:21:A:C8	2.38	0.58
23:DB:2598:A:OP1	25:DC:233:GLY:HA2	2.03	0.58
26:DD:121:THR:C	26:DD:123:LYS:H	2.07	0.58
40:DH:5:LEU:HD13	40:DH:13:GLY:N	2.18	0.58
27:DK:38:ILE:H	27:DK:38:ILE:CD1	2.14	0.58
44:DQ:101:ASP:OD1	44:DQ:104:ALA:HB2	2.02	0.58
44:DQ:40:LYS:HD2	44:DQ:44:TYR:CE1	2.38	0.58
50:DT:16:VAL:H	50:DT:31:VAL:HG23	1.67	0.58
52:DW:18:LYS:HG3	52:DW:19:ARG:CZ	2.33	0.58
52:DW:39:GLN:NE2	52:DW:43:LYS:HB2	2.18	0.58
18:AB:107:ARG:NH2	18:AB:111:LYS:HD3	2.17	0.58
3:AD:48:SER:O	3:AD:52:VAL:HG23	2.03	0.58
6:AG:70:PRO:HA	6:AG:141:HIS:CE1	2.37	0.58
12:AM:3:ILE:HG12	12:AM:56:ARG:NH1	2.18	0.58
23:BB:1021:A:H61	23:BB:1142:A:N6	2.01	0.58
23:BB:1041:G:H2'	23:BB:1042:G:C8	2.38	0.58
23:BB:1097:U:H2'	23:BB:1098:A:H5'	1.85	0.58
23:BB:1548:A:H2'	23:BB:1549:A:H8	1.67	0.58
23:BB:279:A:N6	23:BB:361:G:H1'	2.18	0.58
23:BB:654:A:H2'	23:BB:655:A:H5''	1.86	0.58
23:BB:848:C:H2'	23:BB:849:A:C8	2.38	0.58
25:BC:124:LYS:HE2	25:BC:127:ASN:ND2	2.18	0.58
40:BH:42:LYS:O	40:BH:46:PHE:HB3	2.03	0.58
35:BV:70:ILE:HD13	35:BV:70:ILE:N	2.18	0.58
51:BZ:70:GLU:HA	51:BZ:73:ALA:HB3	1.85	0.58
1:CA:1396:A:H4'	1:CA:1397:C:H5'	1.83	0.58
1:CA:231:U:H2'	1:CA:232:G:H8	1.66	0.58
1:CA:328:C:H4'	1:CA:329:A:C5'	2.33	0.58
1:CA:628:G:H2'	1:CA:629:A:H8	1.68	0.58
4:CE:18:ASN:HB2	4:CE:33:THR:OG1	2.03	0.58
10:CK:88:PRO:HA	10:CK:92:ARG:HD2	1.84	0.58
13:CP:20:VAL:HG23	13:CP:35:ARG:HA	1.85	0.58
15:CR:68:PRO:HB2	15:CR:70:THR:O	2.03	0.58
19:CU:36:PHE:CB	19:CU:40:PRO:HD3	2.28	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1681:G:N3	23:DB:1762:A:H2'	2.17	0.58
23:DB:34:U:H4'	23:DB:34:U:OP2	2.03	0.58
23:DB:2821:A:OP1	26:DD:115:GLY:HA3	2.03	0.58
40:DH:49:ALA:O	40:DH:53:GLU:HB2	2.02	0.58
24:DI:123:ALA:HA	24:DI:126:ARG:HH12	1.67	0.58
27:DK:7:MET:SD	27:DK:20:MET:HB2	2.42	0.58
38:DM:108:VAL:HG21	38:DM:112:LEU:HD12	1.85	0.58
39:DX:8:GLU:HB2	39:DX:13:GLU:OE1	2.02	0.58
1:AA:1226:C:C4	12:AM:102:LYS:HB3	2.38	0.58
1:AA:1229:A:H2'	1:AA:1230:C:H6	1.68	0.58
1:AA:135:C:C2	13:AP:1:MET:HB2	2.38	0.58
1:AA:675:A:H2'	1:AA:676:A:H8	1.68	0.58
1:AA:784:A:H2'	1:AA:785:G:H8	1.68	0.58
1:AA:845:A:N3	1:AA:845:A:H2'	2.18	0.58
2:AC:133:MET:HE1	2:AC:152:VAL:HG21	1.84	0.58
3:AD:32:LYS:O	3:AD:35:GLN:HB2	2.04	0.58
16:AS:39:ILE:HB	16:AS:65:MET:O	2.04	0.58
10:AK:88:PRO:HD3	19:AU:28:LEU:HD11	1.83	0.58
23:BB:1475:G:H4'	23:BB:1476:U:O5'	2.02	0.58
23:BB:1692:U:H2'	23:BB:1694:C:C5	2.38	0.58
23:BB:519:U:H2'	23:BB:520:G:C8	2.38	0.58
23:BB:675:A:H4'	29:BE:62:GLN:HE22	1.68	0.58
23:BB:1790:C:O2'	25:BC:207:ALA:HB2	2.02	0.58
26:BD:106:LYS:O	26:BD:107:VAL:HB	2.03	0.58
48:BG:148:ARG:HD3	48:BG:152:ARG:NH1	2.18	0.58
48:BG:26:LYS:HA	48:BG:32:LEU:N	2.18	0.58
37:BL:73:ILE:HD12	37:BL:106:GLU:HB2	1.84	0.58
38:BM:50:ARG:O	38:BM:53:MET:HB3	2.03	0.58
42:BN:34:ILE:O	42:BN:112:TYR:HA	2.04	0.58
43:BO:88:LYS:HG2	43:BO:116:GLN:HE22	1.67	0.58
28:BP:20:ARG:HB3	28:BP:23:ASP:CG	2.24	0.58
28:BP:31:VAL:HG12	28:BP:38:ARG:O	2.04	0.58
44:BQ:90:ASP:H	49:BR:39:LEU:CD1	2.16	0.58
50:BT:2:ILE:HG12	50:BT:3:ARG:HD3	1.85	0.58
1:CA:1213:A:O2'	1:CA:1214:C:H5'	2.04	0.58
1:CA:1273:C:H2'	1:CA:1274:A:O4'	2.02	0.58
1:CA:468:A:H3'	1:CA:469:C:C6	2.38	0.58
1:CA:956:U:O2'	1:CA:957:U:H5'	2.03	0.58
18:CB:187:ASP:HB3	18:CB:201:GLY:O	2.03	0.58
3:CD:96:ARG:NH1	3:CD:133:SER:HA	2.18	0.58
3:CD:32:LYS:O	3:CD:35:GLN:HB2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CG:145:GLU:HA	6:CG:148:LYS:CB	2.32	0.58
7:CH:31:LEU:HG	7:CH:35:ILE:HD11	1.84	0.58
21:CN:56:PRO:HG2	21:CN:57:SER:H	1.68	0.58
23:DB:1347:A:H2'	23:DB:1348:C:O4'	2.04	0.58
23:DB:1656:C:H2'	23:DB:1657:U:H6	1.68	0.58
23:DB:635:C:O2'	23:DB:639:U:H5''	2.03	0.58
23:DB:759:G:H2'	23:DB:760:G:C8	2.38	0.58
24:DI:32:VAL:HG22	24:DI:60:VAL:HG21	1.86	0.58
41:DJ:96:ARG:O	41:DJ:99:ARG:HG3	2.02	0.58
28:DP:13:LYS:HD2	28:DP:76:HIS:HA	1.84	0.58
27:DK:108:ARG:HH12	28:DP:34:GLY:HA2	1.68	0.58
1:AA:543:U:H2'	1:AA:544:G:C8	2.39	0.58
1:AA:806:C:H2'	1:AA:807:A:C8	2.38	0.58
1:AA:865:A:H2'	1:AA:866:C:C6	2.38	0.58
2:AC:10:ARG:NE	2:AC:181:ILE:HD13	2.19	0.58
3:AD:54:LEU:HA	3:AD:202:LEU:HD22	1.85	0.58
1:AA:921:U:O2	4:AE:23:THR:HG23	2.03	0.58
5:AF:66:ALA:HB1	5:AF:67:PRO:HD2	1.85	0.58
6:AG:149:ALA:H	10:AK:55:ARG:NH2	2.02	0.58
22:BA:32:U:C4'	22:BA:52:A:H62	2.13	0.58
22:BA:75:G:H2'	22:BA:76:G:H8	1.67	0.58
23:BB:1082:U:C2	23:BB:1086:A:C6	2.91	0.58
48:BG:95:ALA:HB2	48:BG:130:ILE:HD11	1.85	0.58
40:BH:90:LEU:HD21	40:BH:146:VAL:HG11	1.84	0.58
27:BK:40:LYS:NZ	27:BK:59:LYS:HE3	2.19	0.58
38:BM:61:GLY:HA2	38:BM:107:GLY:HA3	1.85	0.58
1:CA:1342:C:H2'	1:CA:1343:G:H8	1.69	0.58
1:CA:1370:G:O2'	1:CA:1371:G:H5'	2.04	0.58
1:CA:449:G:H2'	1:CA:450:G:H8	1.67	0.58
1:CA:738:C:H2'	1:CA:739:C:C6	2.37	0.58
1:CA:820:U:H4'	1:CA:821:G:OP2	2.04	0.58
1:CA:865:A:H2	1:CA:918:A:H4'	1.69	0.58
4:CE:136:VAL:HG13	4:CE:137:ARG:H	1.69	0.58
4:CE:95:MET:HA	4:CE:124:ALA:HB2	1.84	0.58
33:D1:52:LYS:HB2	33:D1:52:LYS:HZ2	1.69	0.58
23:DB:1026:G:H2'	23:DB:1027:A:H8	1.68	0.58
23:DB:1584:U:H2'	23:DB:1584:U:O2	2.04	0.58
23:DB:2196:C:O2'	23:DB:2197:U:H5'	2.04	0.58
23:DB:851:C:H2'	23:DB:852:U:H6	1.68	0.58
26:DD:97:SER:HB3	26:DD:99:GLU:HG3	1.85	0.58
47:DF:177:ARG:HA	47:DF:177:ARG:CZ	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DJ:72:LYS:HB2	41:DJ:89:PHE:HB2	1.85	0.58
35:DV:80:HIS:CD2	35:DV:83:LYS:H	2.22	0.58
1:AA:1200:C:O5'	1:AA:1201:A:H3'	2.04	0.58
1:AA:239:U:C5'	1:AA:239:U:H6	2.17	0.58
1:AA:304:U:H2'	1:AA:305:G:C8	2.38	0.58
1:AA:436:C:O2'	1:AA:437:U:H5'	2.04	0.58
1:AA:820:U:H4'	1:AA:821:G:OP2	2.04	0.58
1:AA:948:C:O2'	1:AA:949:A:H5'	2.04	0.58
6:AG:15:PRO:HG2	6:AG:43:TYR:OH	2.04	0.58
6:AG:50:ALA:O	6:AG:54:GLY:N	2.37	0.58
7:AH:81:GLY:O	14:AQ:35:LYS:HD3	2.02	0.58
9:AJ:82:LYS:O	9:AJ:86:ALA:HB2	2.04	0.58
1:AA:707:U:H4'	10:AK:21:HIS:CD2	2.39	0.58
17:AT:47:GLN:HG2	17:AT:82:ILE:HD13	1.84	0.58
17:AT:53:MET:HA	17:AT:56:ILE:HD12	1.85	0.58
17:AT:70:LYS:O	17:AT:73:ARG:HG2	2.03	0.58
32:B4:22:VAL:O	32:B4:24:ARG:HG3	2.03	0.58
23:BB:1593:A:H2'	23:BB:1594:U:C6	2.38	0.58
25:BC:159:THR:O	25:BC:194:VAL:HG12	2.04	0.58
26:BD:159:LYS:HA	26:BD:159:LYS:HZ3	1.68	0.58
26:BD:4:LEU:HD23	26:BD:101:PHE:CE1	2.39	0.58
48:BG:62:ALA:O	48:BG:66:THR:HG22	2.03	0.58
27:BK:99:ILE:HD13	27:BK:118:LEU:HD22	1.85	0.58
37:BL:93:ASN:O	37:BL:95:LEU:N	2.36	0.58
38:BM:69:PRO:HB2	38:BM:92:TRP:HB3	1.85	0.58
31:B0:53:VAL:HG12	42:BN:118:ARG:HH12	1.68	0.58
50:BT:39:THR:HG22	50:BT:42:GLU:HG2	1.86	0.58
52:BW:41:GLY:HA2	52:BW:44:PHE:CD2	2.38	0.58
1:CA:1027:C:H2'	1:CA:1028:C:H6	1.69	0.58
1:CA:106:C:H2'	1:CA:107:G:O4'	2.04	0.58
18:CB:16:GLY:HA2	18:CB:40:ILE:HD12	1.85	0.58
18:CB:33:ALA:HB2	18:CB:38:HIS:HA	1.85	0.58
2:CC:46:LEU:HB3	2:CC:49:ALA:HB3	1.86	0.58
5:CF:6:ILE:HD12	5:CF:7:VAL:N	2.18	0.58
5:CF:86:ARG:CZ	15:CR:63:TYR:HB3	2.34	0.58
12:CM:15:VAL:HG23	12:CM:16:ILE:H	1.69	0.58
12:CM:17:ALA:HB2	12:CM:44:ILE:HD11	1.86	0.58
20:CO:45:GLU:O	20:CO:47:LYS:N	2.37	0.58
10:CK:124:LYS:O	19:CU:33:ARG:NE	2.36	0.58
23:DB:1041:G:H2'	23:DB:1042:G:H8	1.67	0.58
23:DB:1647:U:H3'	23:DB:1647:U:P	2.43	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1683:U:H2'	23:DB:1684:G:H8	1.69	0.58
23:DB:1725:U:H2'	23:DB:1726:C:H6	1.68	0.58
23:DB:1794:A:H2'	23:DB:1795:C:H6	1.68	0.58
23:DB:479:A:N3	23:DB:481:G:H5''	2.17	0.58
23:DB:532:A:O2'	23:DB:2021:C:H5	1.86	0.58
23:DB:974:G:H1'	23:DB:975:A:H8	1.69	0.58
29:DE:149:ILE:O	29:DE:188:MET:HA	2.04	0.58
47:DF:131:VAL:C	47:DF:133:GLU:H	2.07	0.58
40:DH:2:GLN:HB3	40:DH:18:GLN:CG	2.34	0.58
27:DK:109:SER:HB2	27:DK:111:LYS:HE2	1.84	0.58
42:DN:57:THR:HG22	42:DN:58:ASP:H	1.69	0.58
50:DT:14:PRO:HA	50:DT:32:LEU:HB2	1.84	0.58
39:DX:14:LEU:CD2	39:DX:57:LEU:HD21	2.34	0.58
1:AA:501:C:H2'	1:AA:502:A:C8	2.38	0.58
2:AC:104:GLU:O	2:AC:105:VAL:HG13	2.04	0.58
5:AF:17:GLN:O	5:AF:21:MET:HG3	2.04	0.58
12:AM:2:ARG:HG3	12:AM:6:ILE:N	2.18	0.58
23:BB:1013:C:H2'	23:BB:1014:A:H8	1.69	0.58
23:BB:806:C:O2'	23:BB:807:U:H5'	2.04	0.58
26:BD:117:GLY:HA2	26:BD:164:GLN:NE2	2.19	0.58
29:BE:73:ILE:O	29:BE:73:ILE:HG12	2.04	0.58
40:BH:94:ILE:HG22	40:BH:99:ILE:HG12	1.84	0.58
24:BI:27:LEU:CD2	24:BI:27:LEU:H	2.15	0.58
41:BJ:72:LYS:HB2	41:BJ:89:PHE:HB2	1.86	0.58
43:BO:58:ILE:O	43:BO:62:LEU:HG	2.03	0.58
1:CA:631:C:H3'	1:CA:632:U:H5'	1.85	0.58
1:CA:662:U:H2'	1:CA:663:A:H8	1.69	0.58
1:CA:840:C:C4	1:CA:842:U:H4'	2.39	0.58
2:CC:39:ARG:NE	2:CC:56:ILE:HD11	2.19	0.58
1:CA:537:G:H5''	11:CL:109:ARG:NH1	2.18	0.58
12:CM:17:ALA:CB	12:CM:44:ILE:HD11	2.34	0.58
16:CS:21:ALA:HA	16:CS:24:SER:OG	2.03	0.58
16:CS:58:PRO:O	16:CS:59:VAL:HG13	2.04	0.58
17:CT:61:ALA:HA	17:CT:66:ILE:HB	1.85	0.58
23:DB:1061:U:O4'	23:DB:1070:A:H1'	2.03	0.58
23:DB:1220:G:H2'	23:DB:1221:C:H6	1.69	0.58
23:DB:2256:G:H2'	23:DB:2257:U:C6	2.38	0.58
26:DD:34:VAL:HB	26:DD:48:ILE:HD11	1.84	0.58
47:DF:34:THR:O	47:DF:89:THR:HA	2.04	0.58
37:DL:73:ILE:HD12	37:DL:106:GLU:HB2	1.86	0.58
37:DL:4:ASN:ND2	37:DL:4:ASN:N	2.51	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DP:13:LYS:HG2	28:DP:76:HIS:ND1	2.19	0.58
23:DB:996:A:O3'	44:DQ:91:ARG:HG2	2.04	0.58
52:DW:24:ARG:HD3	52:DW:65:LYS:HE3	1.85	0.58
1:AA:469:C:H2'	1:AA:470:C:H6	1.68	0.58
1:AA:993:G:N3	1:AA:993:G:H2'	2.19	0.58
2:AC:111:ASP:O	2:AC:115:VAL:HG23	2.03	0.58
3:AD:90:LEU:HD21	3:AD:196:GLU:HB3	1.86	0.58
6:AG:53:SER:OG	6:AG:55:LYS:HG2	2.04	0.58
6:AG:78:ARG:HB3	6:AG:78:ARG:NH1	2.18	0.58
8:AI:22:PRO:HA	8:AI:60:LEU:HB2	1.86	0.58
6:AG:149:ALA:HB3	10:AK:55:ARG:NH2	2.18	0.58
23:BB:1326:U:O2'	23:BB:1327:A:H5'	2.04	0.58
23:BB:1729:U:H5''	23:BB:1730:C:O2	2.03	0.58
23:BB:1819:A:OP1	25:BC:154:ALA:HA	2.03	0.58
23:BB:2078:C:O2'	23:BB:2079:U:H5'	2.02	0.58
23:BB:2598:A:OP1	25:BC:233:GLY:HA2	2.03	0.58
23:BB:532:A:O2'	23:BB:2021:C:H5	1.87	0.58
23:BB:901:C:H2'	23:BB:902:C:O4'	2.04	0.58
23:BB:784:G:C6	25:BC:227:VAL:HG11	2.39	0.58
47:BF:2:LYS:HE3	47:BF:2:LYS:H	1.67	0.58
41:BJ:29:ALA:HA	41:BJ:32:LEU:HB2	1.86	0.58
42:BN:87:PHE:HB3	42:BN:90:ARG:HB2	1.85	0.58
43:BO:72:ALA:HB2	43:BO:105:ALA:HB1	1.85	0.58
51:BZ:3:ARG:HG2	51:BZ:33:LEU:HD22	1.85	0.58
1:CA:1181:G:O2'	1:CA:1182:G:N7	2.36	0.58
23:DB:1458:U:H4'	23:DB:1459:G:O5'	2.04	0.58
23:DB:2671:G:H2'	23:DB:2672:U:H6	1.68	0.58
23:DB:2693:G:H2'	23:DB:2694:G:H8	1.69	0.58
23:DB:2788:C:H2'	23:DB:2789:C:C6	2.38	0.58
23:DB:856:G:C1'	52:DW:23:LYS:HB3	2.33	0.58
25:DC:132:ARG:HH22	25:DC:169:ALA:HA	1.68	0.58
26:DD:114:LYS:HD2	26:DD:116:LYS:NZ	2.18	0.58
47:DF:97:GLU:O	47:DF:100:GLU:HB2	2.03	0.58
43:DO:111:ARG:HG2	43:DO:117:PHE:H	1.69	0.58
46:DU:86:PHE:HB3	46:DU:90:LYS:O	2.04	0.58
35:DV:61:LEU:O	35:DV:71:LYS:HA	2.03	0.58
52:DW:49:ASN:CB	52:DW:60:ALA:HA	2.33	0.58
52:DW:46:ALA:HB2	52:DW:78:PHE:HD1	1.68	0.58
1:AA:792:A:H1'	1:AA:794:A:N7	2.19	0.58
1:AA:909:A:H2'	1:AA:910:C:O4'	2.04	0.58
18:AB:22:TRP:HA	18:AB:189:ASN:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:55:VAL:HG12	2:AC:56:ILE:H	1.66	0.58
8:AI:11:ARG:HA	8:AI:105:ARG:HE	1.69	0.58
12:AM:58:GLU:HA	12:AM:61:LYS:HE3	1.86	0.58
31:B0:8:THR:HG23	31:B0:11:LYS:H	1.69	0.58
23:BB:1115:G:H2'	23:BB:1116:G:H8	1.68	0.58
23:BB:1347:A:H2'	23:BB:1348:C:O4'	2.04	0.58
23:BB:160:A:H2'	23:BB:161:A:C8	2.39	0.58
23:BB:1656:C:H2'	23:BB:1657:U:H6	1.68	0.58
40:BH:2:GLN:HB3	40:BH:18:GLN:HG3	1.86	0.58
27:BK:17:ARG:HB2	27:BK:45:GLU:HB3	1.85	0.58
50:BT:59:ASN:O	50:BT:84:TYR:HB2	2.03	0.58
1:CA:1296:C:C4'	1:CA:1302:C:H41	2.17	0.58
6:CG:62:GLU:O	6:CG:66:GLU:HG3	2.04	0.58
11:CL:64:SER:OG	11:CL:96:THR:HG23	2.03	0.58
12:CM:104:ASN:N	12:CM:106:ARG:HH21	2.00	0.58
34:D3:39:ARG:O	34:D3:43:LEU:HG	2.03	0.58
32:D4:27:CYS:SG	32:D4:29:ALA:HB3	2.43	0.58
22:DA:11:C:H5'	52:DW:71:LYS:HE3	1.86	0.58
23:DB:132:G:H2'	23:DB:133:U:C6	2.38	0.58
23:DB:1439:A:N7	23:DB:1440:U:C2	2.72	0.58
23:DB:1536:C:H4'	23:DB:1537:G:C4	2.39	0.58
23:DB:2087:G:H2'	23:DB:2088:A:H8	1.68	0.58
23:DB:2105:U:H2'	23:DB:2106:U:C6	2.39	0.58
23:DB:2229:U:H2'	23:DB:2230:G:C8	2.39	0.58
23:DB:2840:C:H2'	23:DB:2841:C:H6	1.67	0.58
23:DB:633:A:O5'	23:DB:633:A:H8	1.86	0.58
23:DB:801:G:H4'	23:DB:802:A:OP2	2.03	0.58
26:DD:8:LYS:HG2	26:DD:197:THR:H	1.69	0.58
23:DB:1099:G:C8	24:DI:3:LYS:CA	2.82	0.58
41:DJ:58:ASN:N	41:DJ:127:GLY:HA2	2.19	0.58
38:DM:42:THR:O	38:DM:44:ARG:N	2.37	0.58
23:DB:988:A:P	30:DY:11:SER:HB3	2.44	0.58
1:AA:1288:A:H2'	1:AA:1289:A:C8	2.39	0.58
1:AA:1396:A:H4'	1:AA:1397:C:H5'	1.85	0.58
2:AC:188:ALA:O	2:AC:194:VAL:HG13	2.04	0.58
9:AJ:8:ILE:CD1	9:AJ:75:ASP:HA	2.33	0.58
13:AP:78:VAL:O	13:AP:80:LYS:N	2.37	0.58
23:BB:2054:A:H2'	31:B0:4:GLN:OE1	2.04	0.58
23:BB:2312:U:H4'	47:BF:84:ILE:HG21	1.86	0.58
23:BB:2840:C:H2'	23:BB:2841:C:H6	1.69	0.58
23:BB:2898:U:H2'	23:BB:2899:A:H8	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:828:U:H4'	23:BB:831:G:N1	2.18	0.58
25:BC:136:VAL:HG12	25:BC:137:GLY:N	2.19	0.58
25:BC:90:ILE:HD13	25:BC:104:LEU:HA	1.85	0.58
26:BD:121:THR:C	26:BD:123:LYS:H	2.07	0.58
29:BE:21:ARG:HG3	29:BE:22:ASP:N	2.17	0.58
48:BG:84:LYS:CG	48:BG:131:VAL:HB	2.34	0.58
41:BJ:57:LEU:HD23	41:BJ:128:ASN:HA	1.85	0.58
37:BL:95:LEU:HB2	37:BL:101:ILE:HG13	1.85	0.58
42:BN:2:ARG:HB3	42:BN:2:ARG:NH1	2.18	0.58
18:CB:160:LEU:HB3	18:CB:181:PRO:O	2.04	0.58
13:CP:7:ALA:O	13:CP:17:TYR:HA	2.04	0.58
17:CT:53:MET:HA	17:CT:56:ILE:HD12	1.85	0.58
23:DB:2615:U:H1'	31:D0:3:GLN:HB3	1.84	0.58
36:D2:19:ARG:HB3	36:D2:19:ARG:NH2	2.19	0.58
47:DF:102:LEU:HD22	47:DF:106:ALA:CB	2.34	0.58
48:DG:84:LYS:CG	48:DG:131:VAL:HB	2.34	0.58
48:DG:30:GLY:O	48:DG:78:VAL:HG12	2.04	0.58
24:DI:99:LYS:HD3	24:DI:99:LYS:H	1.69	0.58
49:DR:29:THR:O	49:DR:63:VAL:HG23	2.04	0.58
46:DU:3:LYS:HA	46:DU:82:VAL:HG11	1.85	0.58
1:AA:1006:G:H2'	1:AA:1006:G:N3	2.18	0.57
1:AA:105:G:H2'	1:AA:106:C:C6	2.39	0.57
1:AA:738:C:H2'	1:AA:739:C:C6	2.39	0.57
18:AB:131:LYS:HB3	18:AB:131:LYS:NZ	2.18	0.57
3:AD:32:LYS:HA	3:AD:35:GLN:HE21	1.69	0.57
3:AD:78:ALA:HA	3:AD:81:LEU:HD12	1.86	0.57
9:AJ:102:LEU:HD12	9:AJ:102:LEU:H	1.68	0.57
12:AM:28:ARG:HH22	12:AM:61:LYS:HB2	1.69	0.57
1:AA:237:G:H5'	14:AQ:26:ARG:HH21	1.69	0.57
16:AS:39:ILE:HD11	16:AS:73:PHE:CE1	2.38	0.57
23:BB:1170:C:H2'	23:BB:1171:G:C8	2.39	0.57
23:BB:1794:A:H2'	23:BB:1795:C:C6	2.39	0.57
23:BB:1831:G:O2'	23:BB:1832:C:H5'	2.04	0.57
23:BB:2262:U:H2'	23:BB:2263:C:H6	1.69	0.57
23:BB:38:A:N3	29:BE:43:THR:HB	2.18	0.57
23:BB:392:U:O2'	23:BB:393:C:H5'	2.03	0.57
25:BC:121:ALA:HB3	25:BC:129:LEU:HD21	1.86	0.57
26:BD:38:LYS:HD3	26:BD:45:TYR:OH	2.04	0.57
29:BE:58:LYS:HE2	29:BE:60:TRP:CD1	2.32	0.57
47:BF:131:VAL:C	47:BF:133:GLU:H	2.08	0.57
40:BH:5:LEU:HB3	40:BH:13:GLY:HA2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BM:71:LYS:HB3	38:BM:93:VAL:HG12	1.86	0.57
50:BT:57:VAL:HG22	50:BT:58:VAL:N	2.13	0.57
1:CA:409:U:H2'	1:CA:410:G:C8	2.39	0.57
18:CB:31:PHE:N	18:CB:41:ASN:HB2	2.19	0.57
7:CH:45:ILE:HA	7:CH:63:LYS:HB2	1.84	0.57
14:CQ:46:HIS:HB2	14:CQ:70:LYS:HE2	1.86	0.57
23:DB:1251:C:H2'	44:DQ:5:ARG:NH1	2.19	0.57
23:DB:1846:G:O2'	23:DB:1847:A:H5'	2.04	0.57
23:DB:2187:U:H2'	23:DB:2188:U:H6	1.69	0.57
26:DD:38:LYS:HD3	26:DD:45:TYR:OH	2.04	0.57
26:DD:34:VAL:HG12	26:DD:94:GLN:H	1.69	0.57
47:DF:14:LYS:O	47:DF:18:GLU:HB2	2.04	0.57
27:DK:102:PRO:HA	27:DK:120:PRO:HB3	1.85	0.57
37:DL:68:SER:C	37:DL:70:LYS:H	2.07	0.57
45:DS:74:ILE:HD12	45:DS:104:THR:O	2.03	0.57
50:DT:59:ASN:O	50:DT:84:TYR:HB2	2.03	0.57
35:DV:72:VAL:CG1	35:DV:93:ARG:HA	2.33	0.57
23:DB:2330:G:N3	52:DW:38:ARG:O	2.37	0.57
51:DZ:70:GLU:HA	51:DZ:73:ALA:HB3	1.86	0.57
1:AA:1364:U:O2'	1:AA:1365:G:H5'	2.05	0.57
1:AA:235:C:H2'	1:AA:236:A:C8	2.40	0.57
1:AA:847:G:H2'	1:AA:848:C:C6	2.40	0.57
1:AA:85:U:O3'	1:AA:86:G:H4'	2.04	0.57
4:AE:18:ASN:HB2	4:AE:33:THR:OG1	2.04	0.57
6:AG:29:LEU:HD13	6:AG:42:VAL:HG22	1.86	0.57
6:AG:57:GLU:CD	6:AG:57:GLU:H	2.07	0.57
11:AL:113:ARG:HG2	11:AL:118:VAL:HB	1.86	0.57
14:AQ:16:MET:CB	14:AQ:19:SER:HB2	2.34	0.57
16:AS:10:ILE:HG21	16:AS:40:PHE:HZ	1.69	0.57
1:AA:1313:U:H3'	16:AS:5:LYS:HG2	1.85	0.57
36:B2:30:VAL:HA	36:B2:33:ARG:HD3	1.86	0.57
23:BB:1166:G:H2'	23:BB:1167:C:H6	1.69	0.57
23:BB:1640:A:H2'	23:BB:1641:A:C8	2.39	0.57
23:BB:1846:G:O2'	23:BB:1847:A:H5'	2.04	0.57
23:BB:2031:A:H5'	23:BB:2031:A:H8	1.68	0.57
23:BB:2019:A:H2	23:BB:2035:G:H22	1.52	0.57
23:BB:2875:C:H2'	23:BB:2876:G:H8	1.69	0.57
23:BB:942:G:O2'	23:BB:943:A:H5'	2.04	0.57
26:BD:46:ARG:NH1	26:BD:86:GLU:H	2.02	0.57
26:BD:97:SER:HB3	26:BD:99:GLU:HG3	1.86	0.57
23:BB:616:A:H4'	29:BE:101:TYR:OH	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:38:A:O2'	29:BE:43:THR:HA	2.04	0.57
23:BB:955:U:P	38:BM:86:LYS:HZ3	2.27	0.57
28:BP:24:THR:O	28:BP:25:VAL:HG22	2.04	0.57
44:BQ:94:LEU:O	44:BQ:97:ILE:HG23	2.03	0.57
45:BS:15:GLN:O	45:BS:19:LEU:HB2	2.04	0.57
45:BS:82:MET:HE1	45:BS:84:ARG:HH22	1.68	0.57
39:BX:12:GLU:HA	39:BX:15:ASN:HD21	1.69	0.57
39:BX:23:ARG:H	39:BX:26:PHE:HB3	1.68	0.57
1:CA:1458:G:H2'	1:CA:1459:G:H8	1.69	0.57
1:CA:1487:G:O2'	1:CA:1488:G:H5'	2.04	0.57
1:CA:1493:A:C3'	1:CA:1493:A:N3	2.67	0.57
1:CA:501:C:H2'	1:CA:502:A:H8	1.68	0.57
18:CB:22:TRP:CG	18:CB:23:ASN:N	2.69	0.57
2:CC:46:LEU:HD11	2:CC:75:VAL:HG13	1.85	0.57
4:CE:37:VAL:HA	4:CE:47:PHE:HA	1.86	0.57
1:CA:1359:C:H3'	21:CN:74:ARG:HH22	1.68	0.57
33:D1:10:LEU:O	33:D1:19:PHE:HB2	2.03	0.57
23:DB:1042:G:H2'	23:DB:1043:C:C6	2.40	0.57
23:DB:1351:C:H2'	23:DB:1352:U:O4'	2.04	0.57
23:DB:1983:G:H4'	23:DB:2606:C:H4'	1.85	0.57
23:DB:300:A:H3'	46:DU:81:ARG:HH12	1.69	0.57
23:DB:516:C:O2'	23:DB:517:C:H5'	2.04	0.57
23:DB:725:G:H2'	23:DB:726:G:O4'	2.05	0.57
25:DC:103:ILE:HG22	25:DC:105:ALA:H	1.69	0.57
25:DC:77:VAL:HG23	25:DC:112:GLY:H	1.65	0.57
26:DD:178:VAL:HG12	26:DD:179:ARG:HG3	1.87	0.57
48:DG:26:LYS:HA	48:DG:32:LEU:N	2.16	0.57
35:DV:68:LYS:H	35:DV:68:LYS:HD3	1.69	0.57
52:DW:41:GLY:HA2	52:DW:44:PHE:CD2	2.39	0.57
1:AA:1013:G:H2'	1:AA:1015:G:OP2	2.03	0.57
1:AA:960:U:H1'	1:AA:1222:G:O2'	2.04	0.57
1:AA:1338:G:H2'	1:AA:1339:A:C8	2.38	0.57
1:AA:976:G:H22	1:AA:1362:A:H3'	1.68	0.57
1:AA:279:A:C5'	1:AA:280:C:H3'	2.34	0.57
1:AA:33:A:H2'	1:AA:34:C:C6	2.39	0.57
1:AA:45:G:H2'	1:AA:46:G:H8	1.68	0.57
1:AA:843:U:H6	1:AA:843:U:H5''	1.68	0.57
8:AI:114:LYS:H	8:AI:120:ALA:HA	1.68	0.57
1:AA:1129:C:H5''	8:AI:17:ARG:NH2	2.20	0.57
8:AI:40:ARG:CA	8:AI:44:ARG:HD3	2.33	0.57
12:AM:77:LYS:HA	12:AM:80:MET:HB3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AQ:75:VAL:HG23	14:AQ:76:ARG:H	1.68	0.57
36:B2:21:ARG:HD3	36:B2:30:VAL:HG11	1.85	0.57
23:BB:1414:C:H2'	23:BB:1415:U:H6	1.69	0.57
23:BB:2063:C:O2	23:BB:2450:A:N1	2.37	0.57
23:BB:2400:G:O2'	23:BB:2401:U:H5'	2.04	0.57
23:BB:309:A:N3	23:BB:329:G:O2'	2.37	0.57
23:BB:813:U:H2'	23:BB:814:C:H6	1.69	0.57
23:BB:974:G:H1'	23:BB:975:A:H8	1.69	0.57
26:BD:5:VAL:HG23	26:BD:32:ASN:HD21	1.69	0.57
29:BE:1:MET:HB3	29:BE:14:VAL:HG23	1.85	0.57
47:BF:64:PRO:HA	47:BF:88:VAL:CG2	2.34	0.57
41:BJ:98:GLU:HB3	41:BJ:124:VAL:HG21	1.85	0.57
27:BK:71:ARG:NE	27:BK:71:ARG:HA	2.19	0.57
30:BY:7:THR:HG22	30:BY:8:GLN:N	2.18	0.57
1:CA:1499:A:H2'	1:CA:1500:A:C8	2.35	0.57
18:CB:61:SER:HB2	18:CB:62:ARG:HH11	1.68	0.57
2:CC:5:HIS:CD2	21:CN:88:MET:HB3	2.38	0.57
3:CD:22:SER:N	3:CD:109:THR:HG22	2.12	0.57
16:CS:1:PRO:O	16:CS:2:ARG:HB2	2.03	0.57
23:DB:1080:A:H2'	23:DB:1081:U:C6	2.38	0.57
23:DB:1315:C:H2'	23:DB:1316:U:C6	2.40	0.57
23:DB:1353:A:H2'	23:DB:1354:A:C8	2.39	0.57
23:DB:1599:U:H2'	23:DB:1600:C:C6	2.39	0.57
23:DB:1656:C:H2'	23:DB:1657:U:C6	2.39	0.57
23:DB:1724:G:H2'	23:DB:1725:U:H6	1.69	0.57
23:DB:2849:U:H4'	23:DB:2850:A:H5'	1.86	0.57
23:DB:2861:U:H2'	23:DB:2862:G:C8	2.40	0.57
23:DB:771:G:O2'	23:DB:772:C:H5'	2.04	0.57
23:DB:82:U:H2'	23:DB:83:A:C8	2.39	0.57
26:DD:31:ALA:HA	26:DD:97:SER:HA	1.87	0.57
29:DE:145:ASP:N	29:DE:166:LYS:HB3	2.18	0.57
29:DE:178:VAL:O	29:DE:181:ILE:HG22	2.04	0.57
48:DG:34:ARG:HH11	48:DG:34:ARG:H	1.51	0.57
48:DG:97:VAL:HB	48:DG:124:CYS:HB2	1.86	0.57
42:DN:54:LEU:HD11	42:DN:62:ASN:HB3	1.85	0.57
28:DP:8:GLU:HB3	28:DP:54:LEU:HB2	1.86	0.57
22:DA:98:G:N1	35:DV:14:LYS:HB2	2.13	0.57
1:AA:952:U:H2'	1:AA:953:G:C8	2.39	0.57
6:AG:14:ASP:HB3	6:AG:19:SER:N	2.12	0.57
1:AA:537:G:H5''	11:AL:109:ARG:NH1	2.20	0.57
20:AO:53:ARG:HD2	23:BB:715:A:N6	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AP:40:ASN:HD21	13:AP:43:ALA:N	2.03	0.57
19:AU:20:ARG:H	19:AU:20:ARG:HD2	1.69	0.57
23:BB:1312:U:H3'	56:BB:3321:HOH:O	2.04	0.57
23:BB:1458:U:H4'	23:BB:1459:G:O5'	2.04	0.57
23:BB:2155:U:H2'	23:BB:2156:G:O4'	2.04	0.57
23:BB:2341:G:H2'	23:BB:2342:C:H6	1.69	0.57
23:BB:39:G:H2'	23:BB:40:U:H6	1.68	0.57
23:BB:224:U:O4	23:BB:420:C:H5'	2.05	0.57
23:BB:551:G:H2'	23:BB:552:U:H6	1.69	0.57
23:BB:586:A:H5'	29:BE:84:THR:OG1	2.04	0.57
48:BG:34:ARG:H	48:BG:34:ARG:HH11	1.52	0.57
27:BK:64:ARG:O	27:BK:82:ASN:HA	2.04	0.57
1:CA:285:C:H2'	1:CA:286:C:H6	1.69	0.57
18:CB:44:LYS:O	18:CB:47:PRO:HD2	2.05	0.57
2:CC:13:ILE:HD13	2:CC:13:ILE:N	2.19	0.57
8:CI:93:LEU:HD13	8:CI:97:LEU:HD11	1.87	0.57
10:CK:70:ALA:C	10:CK:72:ALA:H	2.07	0.57
16:CS:15:LEU:HA	16:CS:18:VAL:CG1	2.35	0.57
23:DB:1050:A:H2'	23:DB:1051:G:O4'	2.04	0.57
23:DB:2085:U:O2'	23:DB:2086:U:H5'	2.04	0.57
23:DB:634:C:H2'	23:DB:635:C:C6	2.38	0.57
23:DB:970:U:H2'	23:DB:971:G:C8	2.40	0.57
29:DE:48:THR:HG22	29:DE:86:ALA:HB3	1.86	0.57
47:DF:139:GLU:HG2	47:DF:140:ILE:H	1.68	0.57
24:DI:125:THR:O	24:DI:129:GLU:HG3	2.05	0.57
27:DK:85:VAL:HG21	27:DK:115:ILE:CD1	2.35	0.57
43:DO:28:VAL:HG11	43:DO:92:PHE:CZ	2.39	0.57
52:DW:37:VAL:HB	52:DW:38:ARG:HD3	1.86	0.57
1:AA:1020:G:H2'	1:AA:1021:A:H5'	1.86	0.57
1:AA:1173:U:H2'	1:AA:1174:G:C8	2.39	0.57
1:AA:1315:U:H3'	1:AA:1316:G:C8	2.40	0.57
1:AA:1448:C:H2'	1:AA:1449:C:C6	2.40	0.57
1:AA:279:A:H5'	1:AA:281:G:H5'	1.85	0.57
18:AB:68:PHE:CE1	18:AB:88:GLN:HB3	2.39	0.57
18:AB:67:LEU:HD22	18:AB:89:PHE:O	2.03	0.57
2:AC:66:THR:HG23	2:AC:101:ASN:ND2	2.19	0.57
9:AJ:6:ILE:HD12	9:AJ:76:ILE:HG13	1.87	0.57
13:AP:7:ALA:O	13:AP:17:TYR:HA	2.04	0.57
16:AS:27:LYS:HZ2	16:AS:27:LYS:HB3	1.67	0.57
17:AT:47:GLN:HE21	17:AT:82:ILE:HD11	1.70	0.57
23:BB:1082:U:O4	23:BB:1086:A:C2	2.57	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1687:G:O2'	23:BB:1688:U:H5'	2.04	0.57
23:BB:2134:A:H62	23:BB:2157:G:N2	1.94	0.57
23:BB:1999:C:H5''	23:BB:2723:C:O2'	2.03	0.57
23:BB:2828:G:O2'	23:BB:2829:A:H5'	2.05	0.57
23:BB:34:U:H4'	23:BB:34:U:OP2	2.05	0.57
23:BB:419:U:H2'	23:BB:420:C:C6	2.39	0.57
23:BB:41:C:H2'	23:BB:42:A:O4'	2.04	0.57
23:BB:516:C:O2'	23:BB:517:C:H5'	2.03	0.57
23:BB:617:G:H5''	29:BE:102:ARG:HH11	1.70	0.57
47:BF:134:GLN:NE2	47:BF:136:ILE:HD13	2.19	0.57
47:BF:22:ASN:HD22	47:BF:22:ASN:H	1.52	0.57
40:BH:80:ILE:HD12	40:BH:98:ASP:O	2.04	0.57
38:BM:21:ALA:HB1	38:BM:100:LYS:HE2	1.86	0.57
28:BP:94:ALA:C	28:BP:95:LYS:HD2	2.24	0.57
46:BU:3:LYS:HA	46:BU:82:VAL:HG11	1.85	0.57
23:BB:102:U:C4	39:BX:2:LYS:HD3	2.40	0.57
39:BX:14:LEU:HD22	39:BX:57:LEU:HD21	1.85	0.57
1:CA:335:C:H2'	1:CA:336:A:C8	2.39	0.57
1:CA:33:A:H2'	1:CA:34:C:C6	2.39	0.57
18:CB:107:ARG:HG3	18:CB:108:GLN:HE21	1.69	0.57
18:CB:113:LEU:HD23	18:CB:114:LYS:N	2.20	0.57
2:CC:50:SER:HB2	2:CC:70:ALA:HB3	1.86	0.57
2:CC:20:THR:HG23	2:CC:57:GLU:HG2	1.85	0.57
4:CE:131:ASN:HD21	4:CE:133:ILE:HB	1.69	0.57
6:CG:42:VAL:O	6:CG:46:LEU:HB2	2.05	0.57
8:CI:54:VAL:HG11	8:CI:86:LEU:CD2	2.34	0.57
12:CM:43:LYS:HD2	12:CM:43:LYS:H	1.69	0.57
22:DA:75:G:H1	22:DA:102:G:N2	2.02	0.57
23:DB:1319:C:O2'	23:DB:1320:C:H5'	2.04	0.57
23:DB:2662:A:H2'	23:DB:2663:G:O4'	2.03	0.57
23:DB:286:U:H2'	23:DB:287:G:H8	1.69	0.57
23:DB:302:C:H2'	23:DB:303:G:H8	1.69	0.57
23:DB:802:A:H2'	23:DB:803:U:C6	2.39	0.57
23:DB:832:U:H2'	23:DB:833:A:H8	1.65	0.57
25:DC:92:LEU:HD12	25:DC:93:VAL:N	2.18	0.57
26:DD:33:ARG:HH11	26:DD:51:THR:HG22	1.68	0.57
29:DE:148:ILE:HA	29:DE:187:VAL:HB	1.86	0.57
40:DH:125:THR:HA	40:DH:146:VAL:CB	2.28	0.57
24:DI:121:ILE:CD1	24:DI:121:ILE:H	2.14	0.57
27:DK:2:ILE:HD12	27:DK:2:ILE:N	2.20	0.57
27:DK:41:ILE:HG13	27:DK:42:THR:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DO:88:LYS:HG2	43:DO:116:GLN:HE22	1.68	0.57
44:DQ:111:LYS:HE2	49:DR:50:GLY:HA3	1.86	0.57
1:AA:1050:G:O2'	1:AA:1051:C:H5'	2.05	0.57
1:AA:1195:C:H2'	1:AA:1197:A:O4'	2.05	0.57
1:AA:1343:G:H2'	1:AA:1344:C:H6	1.69	0.57
1:AA:1372:U:H2'	1:AA:1373:G:O4'	2.04	0.57
18:AB:114:LYS:O	18:AB:118:THR:HG23	2.04	0.57
18:AB:86:CYS:HB3	18:AB:88:GLN:CD	2.24	0.57
2:AC:101:ASN:N	2:AC:101:ASN:OD1	2.37	0.57
5:AF:71:ILE:O	5:AF:75:GLU:HG3	2.03	0.57
8:AI:19:PHE:HB2	8:AI:63:TYR:CB	2.33	0.57
12:AM:84:CYS:HA	16:AS:72:GLU:O	2.04	0.57
23:BB:1026:G:H2'	23:BB:1027:A:H8	1.69	0.57
23:BB:1046:A:H3'	23:BB:1047:G:H5''	1.85	0.57
23:BB:1537:G:H3'	23:BB:1537:G:N3	2.19	0.57
23:BB:1656:C:H2'	23:BB:1657:U:C6	2.39	0.57
23:BB:1725:U:H2'	23:BB:1726:C:H6	1.69	0.57
23:BB:2555:U:H2'	23:BB:2556:C:O4'	2.05	0.57
23:BB:2875:C:H2'	23:BB:2876:G:C8	2.40	0.57
23:BB:771:G:O2'	23:BB:772:C:H5'	2.04	0.57
23:BB:876:C:O4'	23:BB:876:C:O2	2.21	0.57
23:BB:743:A:OP1	26:BD:135:GLY:HA2	2.04	0.57
27:BK:24:VAL:HG13	27:BK:33:ALA:HB2	1.86	0.57
38:BM:4:PRO:HG2	38:BM:70:ASP:HA	1.86	0.57
23:BB:2469:A:H4'	38:BM:55:ARG:CZ	2.35	0.57
38:BM:63:ILE:H	38:BM:63:ILE:HD12	1.69	0.57
44:BQ:40:LYS:HD2	44:BQ:44:TYR:CE1	2.39	0.57
49:BR:62:GLU:O	49:BR:96:VAL:HA	2.03	0.57
45:BS:18:ARG:HB3	45:BS:76:VAL:HG22	1.87	0.57
51:BZ:68:LEU:O	51:BZ:72:ARG:HG2	2.04	0.57
1:CA:1081:A:O2'	1:CA:1082:A:H5'	2.04	0.57
1:CA:1396:A:C3'	1:CA:1397:C:H5'	2.34	0.57
1:CA:1508:A:H2'	1:CA:1509:C:C6	2.40	0.57
2:CC:155:ARG:N	2:CC:162:ALA:HA	2.11	0.57
2:CC:112:ALA:HA	2:CC:201:ILE:HD11	1.86	0.57
23:DB:1021:A:H2'	23:DB:1023:U:H5''	1.86	0.57
23:DB:2216:G:H2'	23:DB:2217:G:C8	2.38	0.57
23:DB:2730:C:H2'	23:DB:2731:G:H8	1.70	0.57
23:DB:322:A:C5'	23:DB:340:A:H1'	2.21	0.57
23:DB:93:G:H2'	23:DB:94:A:C8	2.40	0.57
26:DD:106:LYS:O	26:DD:107:VAL:HB	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:116:LYS:HB3	26:DD:118:PHE:CE1	2.40	0.57
41:DJ:1:MET:HG2	41:DJ:2:LYS:HG2	1.87	0.57
42:DN:33:ILE:HA	42:DN:114:GLU:HB2	1.84	0.57
28:DP:4:ILE:HG22	28:DP:5:LYS:N	2.17	0.57
35:DV:40:ILE:N	35:DV:40:ILE:HD13	2.19	0.57
52:DW:35:ILE:HG13	52:DW:57:THR:OG1	2.05	0.57
39:DX:48:ARG:O	39:DX:51:ALA:HB3	2.04	0.57
1:AA:1187:G:H21	21:AN:99:SER:CB	2.15	0.57
1:AA:1432:G:H1'	1:AA:1468:A:N6	2.19	0.57
1:AA:1492:A:C8	1:AA:1493:A:H1'	2.39	0.57
1:AA:335:C:H2'	1:AA:336:A:H8	1.69	0.57
1:AA:449:G:H2'	1:AA:450:G:H8	1.68	0.57
1:AA:557:G:H2'	1:AA:558:G:O4'	2.03	0.57
8:AI:105:ARG:NH1	8:AI:107:ALA:HA	2.19	0.57
20:AO:60:VAL:HG11	23:BB:715:A:H5'	1.86	0.57
15:AR:68:PRO:HB2	15:AR:70:THR:O	2.04	0.57
19:AU:36:PHE:HA	19:AU:39:LYS:HD2	1.85	0.57
23:BB:705:A:H2'	23:BB:706:A:H8	1.68	0.57
29:BE:143:LEU:HB3	29:BE:146:VAL:HG21	1.85	0.57
40:BH:115:VAL:HB	40:BH:130:VAL:CG1	2.30	0.57
27:BK:107:LEU:HD12	27:BK:107:LEU:N	2.18	0.57
38:BM:42:THR:O	38:BM:44:ARG:N	2.37	0.57
51:BZ:32:ASN:HB2	51:BZ:53:ALA:H	1.68	0.57
1:CA:105:G:H2'	1:CA:106:C:C6	2.40	0.57
1:CA:796:C:O2'	1:CA:797:C:H5'	2.04	0.57
18:CB:46:VAL:HA	18:CB:49:PHE:CD2	2.39	0.57
3:CD:54:LEU:HA	3:CD:202:LEU:HD22	1.86	0.57
3:CD:56:GLU:O	3:CD:60:VAL:HG12	2.05	0.57
4:CE:136:VAL:HG13	4:CE:137:ARG:N	2.20	0.57
1:CA:1240:U:OP1	6:CG:115:MET:HB2	2.05	0.57
17:CT:48:LYS:HG3	17:CT:49:ALA:N	2.19	0.57
33:D1:36:LYS:HB2	33:D1:47:ILE:HA	1.87	0.57
23:DB:101:A:OP1	23:DB:101:A:H4'	2.05	0.57
23:DB:1082:U:O4	23:DB:1086:A:C2	2.57	0.57
23:DB:1727:C:H2'	23:DB:1728:C:C6	2.40	0.57
23:DB:1874:C:H2'	23:DB:1875:G:O4'	2.04	0.57
23:DB:2438:U:O2'	23:DB:2439:A:H5''	2.04	0.57
23:DB:2462:C:H2'	23:DB:2463:C:C6	2.40	0.57
23:DB:901:C:H2'	23:DB:902:C:O4'	2.04	0.57
23:DB:997:G:O2'	23:DB:998:C:H5'	2.05	0.57
47:DF:22:ASN:H	47:DF:22:ASN:HD22	1.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:DG:148:ARG:CD	48:DG:152:ARG:HH11	2.14	0.57
41:DJ:54:ILE:HD12	41:DJ:55:ILE:N	2.20	0.57
38:DM:69:PRO:HB2	38:DM:92:TRP:HB3	1.86	0.57
44:DQ:91:ARG:HB3	44:DQ:93:ILE:HG22	1.85	0.57
44:DQ:94:LEU:O	44:DQ:97:ILE:HG23	2.04	0.57
51:DZ:68:LEU:O	51:DZ:72:ARG:HG2	2.04	0.57
1:AA:1319:A:H4'	1:AA:1320:C:OP1	2.05	0.57
1:AA:1406:U:H2'	1:AA:1407:C:C5'	2.33	0.57
1:AA:195:A:H2'	1:AA:196:A:C8	2.39	0.57
1:AA:598:U:H2'	1:AA:599:C:C6	2.39	0.57
1:AA:865:A:C2	1:AA:918:A:H4'	2.39	0.57
2:AC:149:LYS:O	2:AC:200:TRP:HE3	1.87	0.57
2:AC:91:ALA:HB2	2:AC:98:ALA:HB3	1.86	0.57
20:AO:37:ASN:HA	20:AO:40:GLN:HE22	1.70	0.57
23:BB:2324:U:H1'	23:BB:2337:G:H5'	1.85	0.57
23:BB:2769:U:H2'	23:BB:2770:G:H8	1.69	0.57
23:BB:27:G:HO2'	23:BB:28:A:H8	1.52	0.57
23:BB:634:C:H2'	23:BB:635:C:C6	2.38	0.57
23:BB:742:A:H2'	23:BB:743:A:H8	1.70	0.57
23:BB:997:G:O2'	23:BB:998:C:H5'	2.05	0.57
26:BD:119:ALA:HB1	26:BD:163:GLY:N	2.20	0.57
47:BF:65:LEU:N	47:BF:88:VAL:HG22	2.20	0.57
28:BP:19:PHE:HB2	28:BP:50:ARG:HH12	1.69	0.57
1:CA:1142:G:C2	1:CA:1143:G:H1'	2.39	0.57
1:CA:1488:G:O2'	1:CA:1489:G:H5'	2.05	0.57
18:CB:56:LEU:HB3	18:CB:220:VAL:CG2	2.35	0.57
6:CG:109:LYS:HE2	6:CG:109:LYS:HA	1.87	0.57
13:CP:28:ARG:HH11	13:CP:29:ASN:HD22	1.52	0.57
31:D0:8:THR:HG23	31:D0:11:LYS:H	1.69	0.57
34:D3:20:GLY:HA3	34:D3:48:MET:HE1	1.87	0.57
23:DB:1150:C:H2'	23:DB:1151:A:H8	1.70	0.57
23:DB:1203:U:H3'	23:DB:1204:A:C5'	2.34	0.57
23:DB:1479:G:O2'	23:DB:1480:C:H5'	2.04	0.57
23:DB:1560:G:H2'	23:DB:1561:C:H6	1.68	0.57
23:DB:1571:A:H2'	23:DB:1572:A:C8	2.40	0.57
23:DB:1754:A:H2'	23:DB:1755:A:C8	2.40	0.57
23:DB:2324:U:H1'	23:DB:2337:G:H5'	1.87	0.57
23:DB:2547:A:H5'	23:DB:2566:A:C2	2.39	0.57
23:DB:1999:C:H5''	23:DB:2723:C:O2'	2.05	0.57
23:DB:2728:U:H2'	23:DB:2729:G:H8	1.68	0.57
23:DB:2852:G:H2'	23:DB:2853:C:C6	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:117:GLY:HA2	26:DD:164:GLN:NE2	2.19	0.57
38:DM:4:PRO:HG2	38:DM:70:ASP:HA	1.87	0.57
42:DN:34:ILE:HG22	42:DN:35:LYS:N	2.19	0.57
28:DP:60:VAL:O	28:DP:70:GLU:HA	2.05	0.57
28:DP:94:ALA:C	28:DP:95:LYS:HD2	2.24	0.57
49:DR:14:VAL:HG22	49:DR:15:SER:N	2.20	0.57
1:AA:205:A:H2'	1:AA:206:C:H6	1.70	0.57
18:AB:172:ILE:HG22	18:AB:176:ASN:HD21	1.70	0.57
2:AC:131:ARG:HB3	2:AC:135:ARG:CZ	2.35	0.57
10:AK:70:ALA:C	10:AK:72:ALA:H	2.08	0.57
12:AM:33:LEU:HB3	12:AM:38:ILE:O	2.05	0.57
23:BB:2438:U:O2'	23:BB:2439:A:H5''	2.04	0.57
23:BB:84:A:OP1	46:BU:2:ALA:HB3	2.05	0.57
23:BB:8:C:O2'	23:BB:9:G:H5'	2.05	0.57
23:BB:919:U:H2'	23:BB:920:A:H8	1.69	0.57
26:BD:33:ARG:HD3	26:BD:74:GLU:OE1	2.05	0.57
26:BD:34:VAL:HB	26:BD:48:ILE:HD11	1.85	0.57
47:BF:14:LYS:O	47:BF:18:GLU:HB2	2.04	0.57
48:BG:51:PHE:CE2	48:BG:68:ARG:HA	2.40	0.57
51:BZ:39:TRP:HB2	51:BZ:46:PHE:CE2	2.40	0.57
5:CF:66:ALA:HB1	5:CF:67:PRO:HD2	1.87	0.57
11:CL:113:ARG:HG2	11:CL:118:VAL:HB	1.86	0.57
1:CA:230:G:H5''	13:CP:31:ARG:NH2	2.20	0.57
23:DB:1652:A:H62	42:DN:11:ASN:ND2	1.96	0.57
23:DB:2267:A:C8	23:DB:2267:A:C3'	2.77	0.57
23:DB:639:U:H2'	23:DB:640:C:C6	2.39	0.57
23:DB:814:C:O2'	23:DB:815:C:H5'	2.05	0.57
23:DB:784:G:N1	25:DC:227:VAL:HG11	2.20	0.57
26:DD:54:ALA:CA	26:DD:76:GLY:HA2	2.33	0.57
28:DP:20:ARG:HB3	28:DP:23:ASP:CG	2.24	0.57
1:AA:1149:C:H2'	1:AA:1150:A:C8	2.39	0.57
1:AA:1382:C:H2'	1:AA:1383:C:C6	2.40	0.57
1:AA:631:C:H3'	1:AA:632:U:H5'	1.86	0.57
7:AH:95:MET:HB3	7:AH:99:GLY:H	1.68	0.57
20:AO:62:GLN:O	20:AO:66:LEU:HB2	2.05	0.57
20:AO:69:TYR:HA	20:AO:72:ARG:CZ	2.35	0.57
23:BB:1351:C:H2'	23:BB:1352:U:O4'	2.05	0.57
23:BB:136:G:H2'	23:BB:137:U:C1'	2.35	0.57
23:BB:1560:G:H2'	23:BB:1561:C:H6	1.68	0.57
23:BB:1584:U:O2	23:BB:1584:U:H2'	2.05	0.57
23:BB:18:U:H2'	23:BB:19:A:H8	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:675:A:OP1	29:BE:60:TRP:NE1	2.38	0.57
23:BB:817:C:H2'	23:BB:818:G:O4'	2.05	0.57
23:BB:934:U:H2'	23:BB:935:C:C6	2.40	0.57
25:BC:146:LYS:HB2	25:BC:149:LYS:HB2	1.87	0.57
29:BE:200:LEU:O	29:BE:201:ALA:HB3	2.05	0.57
37:BL:110:VAL:HB	37:BL:127:VAL:HG23	1.85	0.57
28:BP:31:VAL:O	28:BP:32:VAL:HB	2.03	0.57
28:BP:47:ILE:HD11	28:BP:59:THR:HG22	1.86	0.57
50:BT:32:LEU:O	50:BT:83:ALA:HB2	2.05	0.57
46:BU:86:PHE:CG	46:BU:87:GLU:N	2.73	0.57
52:BW:46:ALA:HB2	52:BW:78:PHE:HD1	1.70	0.57
52:BW:49:ASN:O	52:BW:50:VAL:HG13	2.04	0.57
39:BX:14:LEU:CD2	39:BX:57:LEU:HD21	2.35	0.57
51:BZ:21:ALA:HB3	51:BZ:23:ASN:ND2	2.20	0.57
1:CA:205:A:H2'	1:CA:206:C:C6	2.39	0.57
1:CA:377:G:H2'	1:CA:378:G:H8	1.69	0.57
1:CA:499:A:H4'	1:CA:500:G:H5'	1.87	0.57
1:CA:624:C:H2'	1:CA:625:U:C6	2.40	0.57
1:CA:797:C:O2'	1:CA:798:U:H5'	2.04	0.57
1:CA:909:A:H2'	1:CA:910:C:O4'	2.04	0.57
2:CC:119:ILE:O	2:CC:123:LEU:HG	2.05	0.57
5:CF:73:GLU:O	5:CF:77:THR:HG23	2.05	0.57
8:CI:12:LYS:H	8:CI:105:ARG:HH12	1.53	0.57
10:CK:59:PRO:HB2	6:CG:149:ALA:HB1	1.86	0.57
12:CM:18:LEU:HD12	12:CM:29:SER:OG	2.05	0.57
1:CA:1308:U:OP2	12:CM:97:ARG:HD3	2.04	0.57
20:CO:81:LEU:HD23	20:CO:82:ILE:N	2.20	0.57
19:CU:42:THR:O	19:CU:46:ARG:HG3	2.05	0.57
34:D3:22:LYS:HG3	34:D3:47:ALA:O	2.05	0.57
23:DB:1039:A:H2'	23:DB:1040:A:C8	2.40	0.57
23:DB:1199:U:H2'	23:DB:1200:C:C6	2.39	0.57
23:DB:1831:G:O2'	23:DB:1832:C:H5'	2.05	0.57
23:DB:2257:U:O2'	23:DB:2258:C:H5'	2.05	0.57
23:DB:519:U:H2'	23:DB:520:G:C8	2.40	0.57
23:DB:877:A:H61	23:DB:898:C:C2'	2.12	0.57
47:DF:87:LYS:CG	47:DF:88:VAL:H	2.18	0.57
41:DJ:59:ALA:HB1	41:DJ:101:ILE:HD11	1.87	0.57
23:DB:584:C:OP2	44:DQ:5:ARG:HD3	2.05	0.57
1:AA:285:C:H2'	1:AA:286:C:H6	1.69	0.56
1:AA:332:G:O2'	1:AA:333:U:H5'	2.04	0.56
1:AA:711:G:O2'	1:AA:712:A:H5'	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:87:MET:HB2	8:AI:94:ARG:HD3	1.87	0.56
9:AJ:40:ILE:H	9:AJ:74:VAL:H	1.51	0.56
23:BB:1897:G:O2'	23:BB:1898:U:H5'	2.05	0.56
23:BB:2052:A:H4'	26:BD:148:GLN:O	2.05	0.56
23:BB:2443:C:H2'	23:BB:2444:G:C8	2.40	0.56
23:BB:394:C:O2'	23:BB:395:U:H5'	2.05	0.56
47:BF:72:SER:HA	47:BF:78:ILE:HG22	1.87	0.56
41:BJ:45:THR:N	41:BJ:46:PRO:HD3	2.20	0.56
42:BN:19:ALA:C	42:BN:21:PHE:H	2.06	0.56
50:BT:9:LYS:NZ	50:BT:9:LYS:HB3	2.20	0.56
1:CA:1218:C:H2'	1:CA:1219:A:C8	2.39	0.56
1:CA:1237:C:H2'	1:CA:1336:C:H5	1.69	0.56
1:CA:838:G:H2'	1:CA:839:C:C6	2.39	0.56
2:CC:149:LYS:HG2	2:CC:200:TRP:HE3	1.69	0.56
10:CK:33:ILE:CB	10:CK:73:VAL:HG11	2.33	0.56
11:CL:17:LYS:HE3	11:CL:17:LYS:N	2.20	0.56
12:CM:13:HIS:CD2	12:CM:13:HIS:H	2.23	0.56
23:DB:1177:G:H2'	23:DB:1178:C:C6	2.40	0.56
23:DB:2229:U:H2'	23:DB:2230:G:H8	1.68	0.56
23:DB:276:U:H2'	23:DB:278:A:N6	2.13	0.56
23:DB:2898:U:H2'	23:DB:2899:A:H8	1.70	0.56
23:DB:598:U:H2'	23:DB:599:A:H8	1.69	0.56
25:DC:245:THR:C	25:DC:247:TRP:H	2.09	0.56
29:DE:143:LEU:HB3	29:DE:146:VAL:HG21	1.85	0.56
47:DF:37:MET:HG3	47:DF:150:GLY:O	2.05	0.56
48:DG:9:VAL:HA	48:DG:48:THR:CB	2.35	0.56
40:DH:57:LYS:O	40:DH:61:VAL:HB	2.04	0.56
41:DJ:45:THR:N	41:DJ:46:PRO:HD3	2.20	0.56
27:DK:103:VAL:HG23	27:DK:122:VAL:O	2.04	0.56
27:DK:86:LEU:HD23	27:DK:86:LEU:H	1.70	0.56
42:DN:34:ILE:O	42:DN:112:TYR:HA	2.05	0.56
28:DP:24:THR:O	28:DP:25:VAL:HG22	2.04	0.56
50:DT:69:ARG:HG2	50:DT:74:ILE:N	2.20	0.56
46:DU:51:LEU:H	46:DU:53:GLN:NE2	2.02	0.56
1:AA:1095:U:H5''	1:AA:1109:C:O2	2.05	0.56
1:AA:1164:G:H2'	1:AA:1165:U:C6	2.40	0.56
1:AA:1208:C:H2'	1:AA:1209:C:O4'	2.05	0.56
1:AA:1308:U:H2'	1:AA:1309:G:C8	2.40	0.56
1:AA:1348:U:C4'	8:AI:121:ARG:HG3	2.35	0.56
1:AA:1394:A:H5'	56:AA:1962:HOH:O	2.05	0.56
1:AA:556:C:O2'	1:AA:557:G:H5'	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:39:ARG:NE	2:AC:56:ILE:HD11	2.19	0.56
12:AM:14:ALA:HB3	12:AM:40:GLU:HA	1.86	0.56
23:BB:154:U:H2'	23:BB:155:A:H8	1.70	0.56
23:BB:1590:A:H2'	23:BB:1591:A:H8	1.70	0.56
23:BB:216:A:H2'	23:BB:217:A:H8	1.70	0.56
23:BB:2297:A:N1	23:BB:2320:U:H4'	2.20	0.56
23:BB:2835:A:N6	23:BB:2878:U:H2'	2.21	0.56
23:BB:402:A:C2'	23:BB:403:U:H5'	2.35	0.56
25:BC:136:VAL:HG12	25:BC:137:GLY:H	1.70	0.56
25:BC:83:ASP:HB2	25:BC:90:ILE:HG12	1.85	0.56
27:BK:3:GLN:HG2	27:BK:4:GLU:N	2.19	0.56
27:BK:71:ARG:HE	27:BK:71:ARG:HA	1.70	0.56
37:BL:68:SER:C	37:BL:70:LYS:H	2.07	0.56
1:CA:577:G:O2'	1:CA:578:C:H5'	2.05	0.56
1:CA:691:G:H1'	1:CA:696:A:N6	2.20	0.56
1:CA:704:A:H2'	1:CA:705:G:O4'	2.05	0.56
18:CB:187:ASP:H	18:CB:190:SER:HB2	1.69	0.56
13:CP:22:ALA:HA	13:CP:33:ILE:HG13	1.88	0.56
16:CS:60:PHE:N	16:CS:60:PHE:HD2	2.04	0.56
1:CA:1320:C:H1'	16:CS:72:GLU:N	2.19	0.56
17:CT:70:LYS:O	17:CT:73:ARG:HG2	2.04	0.56
23:DB:1166:G:H2'	23:DB:1167:C:H6	1.71	0.56
23:DB:2498:C:O2'	23:DB:2499:C:H5'	2.05	0.56
23:DB:2590:A:H2'	23:DB:2591:C:C6	2.41	0.56
23:DB:598:U:H2'	23:DB:599:A:C8	2.41	0.56
23:DB:648:G:H2'	23:DB:649:G:H8	1.68	0.56
25:DC:136:VAL:HG12	25:DC:137:GLY:H	1.70	0.56
25:DC:146:LYS:HB2	25:DC:149:LYS:HB2	1.86	0.56
47:DF:155:ILE:HG22	47:DF:157:THR:H	1.70	0.56
48:DG:137:LYS:O	48:DG:140:ILE:HG13	2.05	0.56
41:DJ:13:ARG:O	41:DJ:52:ASP:HA	2.05	0.56
27:DK:99:ILE:HD13	27:DK:118:LEU:HD22	1.85	0.56
27:DK:39:ILE:O	27:DK:59:LYS:HA	2.05	0.56
28:DP:3:ILE:HG23	28:DP:4:ILE:H	1.70	0.56
28:DP:63:ILE:HG12	28:DP:68:GLY:HA2	1.88	0.56
28:DP:20:ARG:HH22	28:DP:91:VAL:HG22	1.69	0.56
45:DS:18:ARG:HB3	45:DS:76:VAL:CG2	2.35	0.56
45:DS:81:SER:HA	45:DS:99:ARG:HA	1.87	0.56
50:DT:32:LEU:HG	50:DT:83:ALA:HB2	1.88	0.56
46:DU:12:VAL:HA	46:DU:69:VAL:HG12	1.86	0.56
1:AA:1072:G:H21	18:AB:105:THR:HG21	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1229:A:H2'	1:AA:1230:C:C6	2.39	0.56
1:AA:991:U:H3	1:AA:1212:U:H4'	1.69	0.56
2:AC:122:GLN:HG2	2:AC:127:VAL:HG11	1.86	0.56
6:AG:8:GLN:HG2	6:AG:9:ARG:N	2.20	0.56
14:AQ:59:GLU:HB3	14:AQ:75:VAL:CG2	2.35	0.56
23:BB:1439:A:N7	23:BB:1440:U:C2	2.73	0.56
23:BB:1442:U:H2'	23:BB:1443:U:H6	1.69	0.56
23:BB:1842:G:H2'	23:BB:1843:C:H6	1.70	0.56
26:BD:55:LYS:HE2	26:BD:59:ARG:HB2	1.86	0.56
48:BG:94:ARG:HH21	48:BG:104:LEU:HA	1.71	0.56
50:BT:16:VAL:H	50:BT:31:VAL:HG23	1.71	0.56
35:BV:40:ILE:N	35:BV:40:ILE:HD13	2.19	0.56
23:BB:922:C:H1'	52:BW:22:VAL:CG2	2.35	0.56
1:CA:1192:C:H2'	1:CA:1193:G:O4'	2.05	0.56
1:CA:255:G:H2'	1:CA:256:U:C6	2.40	0.56
1:CA:532:A:N6	2:CC:126:ARG:HH21	2.03	0.56
1:CA:821:G:H4'	56:CA:1750:HOH:O	2.03	0.56
5:CF:17:GLN:O	5:CF:21:MET:HG3	2.05	0.56
8:CI:39:GLY:O	8:CI:40:ARG:HB2	2.06	0.56
16:CS:40:PHE:HB2	16:CS:43:MET:SD	2.44	0.56
23:DB:1268:A:H2'	23:DB:1269:A:O4'	2.05	0.56
23:DB:1326:U:O2'	23:DB:1327:A:H5'	2.05	0.56
23:DB:1590:A:H2'	23:DB:1591:A:H8	1.70	0.56
23:DB:2251:G:H2'	23:DB:2252:G:C8	2.40	0.56
23:DB:2784:U:H2'	23:DB:2785:C:C6	2.41	0.56
23:DB:365:U:H2'	23:DB:366:C:C6	2.40	0.56
23:DB:414:C:H2'	23:DB:415:A:H8	1.67	0.56
23:DB:705:A:H2'	23:DB:706:A:H8	1.70	0.56
23:DB:737:C:O2'	23:DB:738:G:H5'	2.06	0.56
25:DC:136:VAL:HG12	25:DC:137:GLY:N	2.20	0.56
25:DC:191:LEU:CD2	25:DC:192:GLY:H	2.17	0.56
22:DA:41:G:O6	47:DF:68:LYS:HD3	2.04	0.56
24:DI:108:ILE:HG22	24:DI:128:ILE:HD13	1.87	0.56
41:DJ:57:LEU:HD23	41:DJ:128:ASN:HA	1.87	0.56
38:DM:71:LYS:HB3	38:DM:93:VAL:HG12	1.86	0.56
40:DH:27:ARG:CZ	51:DZ:60:ASP:HA	2.34	0.56
4:AE:140:ILE:HG22	4:AE:144:GLU:OE1	2.05	0.56
8:AI:29:ILE:HA	8:AI:64:ILE:CB	2.32	0.56
33:B1:9:LYS:N	33:B1:9:LYS:HD3	2.19	0.56
23:BB:1051:G:H2'	23:BB:1052:C:C6	2.41	0.56
23:BB:1346:G:O2'	23:BB:1347:A:H5'	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1809:A:H2'	23:BB:1810:A:C8	2.41	0.56
23:BB:1859:U:H2'	23:BB:1860:G:C8	2.41	0.56
23:BB:1869:G:H3'	23:BB:1870:C:C5'	2.35	0.56
23:BB:2097:A:H2'	23:BB:2098:U:H6	1.69	0.56
23:BB:345:A:H1'	23:BB:346:A:C2	2.39	0.56
23:BB:370:G:O2'	23:BB:423:A:H3'	2.06	0.56
23:BB:670:A:H4'	23:BB:671:C:C5'	2.34	0.56
23:BB:876:C:H41	23:BB:901:C:H42	1.48	0.56
29:BE:146:VAL:O	29:BE:167:VAL:HA	2.06	0.56
29:BE:58:LYS:HB2	29:BE:60:TRP:HD1	1.71	0.56
48:BG:84:LYS:C	48:BG:85:LYS:HG2	2.26	0.56
40:BH:2:GLN:HB3	40:BH:18:GLN:CG	2.35	0.56
41:BJ:18:VAL:O	41:BJ:56:VAL:HA	2.04	0.56
41:BJ:25:LEU:HD22	41:BJ:26:GLY:H	1.70	0.56
41:BJ:28:LEU:HG	41:BJ:29:ALA:N	2.21	0.56
27:BK:87:LEU:HD12	27:BK:93:GLN:H	1.69	0.56
50:BT:14:PRO:HA	50:BT:32:LEU:CB	2.35	0.56
1:CA:1128:C:H4'	1:CA:1148:U:O2	2.04	0.56
1:CA:266:G:O2'	1:CA:267:C:H3'	2.05	0.56
1:CA:462:G:H3'	1:CA:463:U:H6	1.70	0.56
3:CD:29:THR:HB	3:CD:30:LYS:HZ3	1.70	0.56
6:CG:30:MET:HE3	6:CG:33:GLY:HA2	1.86	0.56
9:CJ:8:ILE:HA	9:CJ:100:ILE:CG2	2.35	0.56
16:CS:21:ALA:HA	16:CS:24:SER:HG	1.70	0.56
16:CS:60:PHE:N	16:CS:60:PHE:CD2	2.74	0.56
23:DB:1599:U:H2'	23:DB:1600:C:H6	1.71	0.56
23:DB:1640:A:H2'	23:DB:1641:A:H8	1.69	0.56
23:DB:172:A:O2'	23:DB:173:A:H5'	2.05	0.56
23:DB:2611:C:O2'	23:DB:2612:C:H5'	2.04	0.56
23:DB:545:U:H2'	23:DB:548:G:OP2	2.06	0.56
26:DD:33:ARG:HH11	26:DD:76:GLY:HA3	1.70	0.56
47:DF:102:LEU:HA	47:DF:106:ALA:CB	2.35	0.56
47:DF:33:ILE:HB	47:DF:90:LEU:HB2	1.87	0.56
23:DB:2748:A:H1'	48:DG:66:THR:HB	1.85	0.56
24:DI:37:PHE:CZ	24:DI:58:ILE:HD11	2.40	0.56
27:DK:24:VAL:HG13	27:DK:33:ALA:HB2	1.86	0.56
43:DO:74:VAL:O	43:DO:77:ALA:HB3	2.05	0.56
44:DQ:111:LYS:HE2	49:DR:50:GLY:CA	2.36	0.56
46:DU:42:LYS:HD2	46:DU:59:GLU:HB2	1.87	0.56
46:DU:5:ARG:NH2	46:DU:93:ARG:HD3	2.21	0.56
1:AA:135:C:N3	13:AP:1:MET:HB2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1360:A:H61	1:AA:1361:G:H21	1.53	0.56
1:AA:1486:G:H2'	1:AA:1487:G:C1'	2.36	0.56
1:AA:1504:G:H4'	1:AA:1505:G:O4'	2.06	0.56
1:AA:202:G:H2'	1:AA:203:G:C8	2.39	0.56
1:AA:616:G:H2'	1:AA:616:G:N3	2.21	0.56
1:AA:861:G:H2'	1:AA:862:C:H6	1.70	0.56
1:AA:880:C:H2'	1:AA:881:G:C8	2.41	0.56
1:AA:919:A:O2'	1:AA:920:U:H5'	2.05	0.56
1:AA:987:G:H2'	1:AA:988:G:H8	1.70	0.56
3:AD:22:SER:N	3:AD:109:THR:HG22	2.12	0.56
3:AD:90:LEU:HD22	3:AD:90:LEU:H	1.70	0.56
8:AI:126:PHE:O	8:AI:128:LYS:N	2.38	0.56
12:AM:87:GLY:HA2	12:AM:90:HIS:CD2	2.41	0.56
13:AP:53:ASP:O	13:AP:57:ILE:HG12	2.05	0.56
23:BB:1724:G:H2'	23:BB:1725:U:H6	1.69	0.56
23:BB:1914:C:O2	23:BB:1914:C:H3'	2.05	0.56
23:BB:2078:C:H2'	23:BB:2079:U:C6	2.39	0.56
25:BC:204:LEU:HD23	25:BC:209:ALA:HB3	1.88	0.56
25:BC:20:ASN:ND2	25:BC:23:LEU:HD13	2.21	0.56
26:BD:107:VAL:HA	26:BD:204:LYS:O	2.06	0.56
26:BD:101:PHE:O	26:BD:180:VAL:HG11	2.05	0.56
24:BI:58:ILE:N	24:BI:58:ILE:HD12	2.20	0.56
23:BB:2548:U:H1'	27:BK:23:LYS:NZ	2.19	0.56
38:BM:18:ARG:C	38:BM:38:ARG:HH22	2.09	0.56
42:BN:54:LEU:HD11	42:BN:62:ASN:HB3	1.88	0.56
28:BP:20:ARG:HH22	28:BP:91:VAL:HG22	1.70	0.56
49:BR:37:GLU:O	49:BR:39:LEU:HD23	2.05	0.56
39:BX:8:GLU:HB2	39:BX:13:GLU:OE1	2.06	0.56
1:CA:1223:C:O5'	1:CA:1224:U:H5''	2.05	0.56
1:CA:1396:A:C4'	1:CA:1397:C:H5'	2.35	0.56
1:CA:1524:C:H2'	1:CA:1525:G:C8	2.41	0.56
1:CA:675:A:H2'	1:CA:676:A:C8	2.41	0.56
1:CA:733:G:O2'	1:CA:734:G:H5'	2.05	0.56
3:CD:199:ILE:HG13	3:CD:200:VAL:N	2.21	0.56
8:CI:103:VAL:HG23	8:CI:104:THR:N	2.19	0.56
9:CJ:42:LEU:HG	9:CJ:43:PRO:HD2	1.87	0.56
16:CS:32:THR:HG22	16:CS:33:TRP:N	2.20	0.56
16:CS:39:ILE:HD12	16:CS:68:HIS:H	1.70	0.56
23:DB:362:A:H3'	23:DB:363:G:C8	2.40	0.56
23:DB:519:U:H2'	23:DB:520:G:H8	1.70	0.56
26:DD:186:LEU:HD21	28:DP:3:ILE:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:108:PRO:C	47:DF:110:ILE:H	2.09	0.56
47:DF:64:PRO:HA	47:DF:88:VAL:CG2	2.35	0.56
40:DH:70:GLU:HA	40:DH:73:ASN:HB2	1.87	0.56
23:DB:1098:A:H3'	24:DI:3:LYS:C	2.26	0.56
41:DJ:93:ILE:O	41:DJ:97:PRO:HG3	2.05	0.56
27:DK:40:LYS:HZ2	27:DK:59:LYS:HE3	1.70	0.56
23:DB:1952:A:OP1	27:DK:42:THR:HG21	2.06	0.56
37:DL:3:LEU:O	37:DL:5:THR:HG23	2.05	0.56
1:CA:1463:U:OP1	28:DP:108:ARG:HD2	2.05	0.56
44:DQ:10:ARG:HB2	44:DQ:10:ARG:CZ	2.34	0.56
23:DB:584:C:P	44:DQ:5:ARG:HD3	2.46	0.56
46:DU:85:ARG:HD3	46:DU:86:PHE:N	2.20	0.56
1:AA:1488:G:O2'	1:AA:1489:G:H5'	2.05	0.56
1:AA:513:C:H2'	1:AA:514:C:C6	2.39	0.56
1:AA:838:G:H2'	1:AA:839:C:C6	2.40	0.56
1:AA:927:G:O2'	1:AA:928:G:H5'	2.06	0.56
9:AJ:12:ALA:HB3	9:AJ:18:ILE:HB	1.87	0.56
10:AK:28:ASN:CG	10:AK:56:LYS:HE3	2.26	0.56
11:AL:14:LYS:HZ3	11:AL:16:ALA:HB3	1.69	0.56
22:BA:59:A:H2'	22:BA:60:C:C6	2.40	0.56
23:BB:1061:U:O4'	23:BB:1070:A:H1'	2.06	0.56
23:BB:2015:A:C2	31:B0:2:VAL:HG22	2.40	0.56
23:BB:2292:U:H2'	23:BB:2293:G:C8	2.41	0.56
23:BB:2384:U:H5''	23:BB:2386:A:OP1	2.05	0.56
23:BB:2393:U:H2'	23:BB:2394:C:O4'	2.06	0.56
23:BB:264:C:C2'	23:BB:265:A:H5''	2.34	0.56
23:BB:2845:U:O3'	28:BP:52:ARG:HD3	2.04	0.56
23:BB:417:C:H2'	23:BB:418:C:H6	1.71	0.56
23:BB:2531:A:H5'	48:BG:156:TYR:CE2	2.41	0.56
27:BK:85:VAL:HG21	27:BK:115:ILE:HD11	1.86	0.56
37:BL:4:ASN:N	37:BL:4:ASN:ND2	2.54	0.56
44:BQ:77:LYS:HA	44:BQ:80:ASN:HB3	1.87	0.56
46:BU:32:LYS:HA	46:BU:65:GLN:HA	1.86	0.56
46:BU:82:VAL:H	46:BU:96:LYS:NZ	2.04	0.56
1:CA:1221:G:H4'	16:CS:76:THR:CG2	2.28	0.56
1:CA:1458:G:H4'	17:CT:22:SER:HB2	1.86	0.56
1:CA:285:C:H2'	1:CA:286:C:C6	2.41	0.56
18:CB:116:LEU:HD11	18:CB:139:GLU:OE1	2.06	0.56
18:CB:67:LEU:HD22	18:CB:91:VAL:HG23	1.88	0.56
6:CG:100:MET:O	6:CG:104:VAL:HG23	2.06	0.56
7:CH:108:GLY:O	7:CH:110:MET:HG3	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:6:ILE:O	9:CJ:8:ILE:HD12	2.05	0.56
19:CU:19:LYS:HD2	19:CU:19:LYS:N	2.20	0.56
23:DB:1552:A:H2'	23:DB:1553:A:H5'	1.87	0.56
23:DB:2231:U:H2'	23:DB:2232:C:C6	2.40	0.56
23:DB:224:U:O4	23:DB:420:C:H5'	2.06	0.56
23:DB:2330:G:H1'	52:DW:38:ARG:CB	2.34	0.56
26:DD:107:VAL:HA	26:DD:204:LYS:O	2.05	0.56
26:DD:56:LYS:HD2	26:DD:59:ARG:HG3	1.86	0.56
29:DE:21:ARG:HG3	29:DE:22:ASP:O	2.05	0.56
47:DF:78:ILE:N	47:DF:78:ILE:HD12	2.20	0.56
28:DP:89:GLY:HA2	28:DP:111:GLU:C	2.26	0.56
28:DP:21:PRO:CG	28:DP:91:VAL:HG21	2.34	0.56
51:DZ:59:ILE:HD13	51:DZ:67:VAL:HG21	1.88	0.56
1:AA:195:A:H1'	1:AA:222:C:O2'	2.06	0.56
1:AA:451:A:H5'	13:AP:70:ARG:HH22	1.70	0.56
1:AA:484:G:H5'	1:AA:486:U:H5'	1.88	0.56
1:AA:678:U:H2'	1:AA:679:C:C6	2.40	0.56
3:AD:3:TYR:C	3:AD:4:LEU:HD12	2.26	0.56
8:AI:49:GLN:N	8:AI:50:PRO:HD2	2.21	0.56
1:AA:255:G:H4'	14:AQ:18:LYS:HD2	1.87	0.56
16:AS:39:ILE:HD13	16:AS:65:MET:HB3	1.87	0.56
23:BB:1635:A:C2'	23:BB:1636:U:H5'	2.36	0.56
23:BB:1704:C:H2'	23:BB:1705:A:H8	1.71	0.56
23:BB:2748:A:H4'	48:BG:3:VAL:HG21	1.88	0.56
23:BB:970:U:H2'	23:BB:971:G:C8	2.41	0.56
5:AF:80:PHE:CE1	25:BC:123:ILE:HG12	2.41	0.56
25:BC:202:ARG:NH1	25:BC:204:LEU:HD11	2.20	0.56
26:BD:109:VAL:HG11	26:BD:193:VAL:CG1	2.35	0.56
26:BD:116:LYS:HB3	26:BD:118:PHE:CE1	2.41	0.56
29:BE:118:LEU:HD21	29:BE:188:MET:HE3	1.88	0.56
47:BF:78:ILE:HD12	47:BF:78:ILE:N	2.21	0.56
27:BK:108:ARG:HH12	28:BP:34:GLY:HA2	1.70	0.56
42:BN:85:PRO:HA	42:BN:88:ALA:HB2	1.86	0.56
43:BO:11:ALA:HB2	43:BO:96:GLY:N	2.20	0.56
51:BZ:32:ASN:C	51:BZ:33:LEU:HD12	2.26	0.56
1:CA:1027:C:H2'	1:CA:1028:C:C6	2.41	0.56
1:CA:114:U:O2'	1:CA:115:G:H5'	2.06	0.56
1:CA:1289:A:H3'	1:CA:1290:G:H8	1.70	0.56
1:CA:1513:A:H2'	1:CA:1514:G:C8	2.40	0.56
1:CA:213:G:H3'	1:CA:214:C:H6	1.71	0.56
18:CB:19:THR:HG23	18:CB:20:ARG:H	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:146:LYS:HB2	2:CC:202:PHE:CD2	2.41	0.56
9:CJ:65:TYR:HA	21:CN:98:ALA:H	1.71	0.56
20:CO:88:ARG:O	20:CO:88:ARG:HG3	2.05	0.56
23:DB:1309:G:OP1	36:D2:9:VAL:HG12	2.05	0.56
34:D3:18:LYS:HD2	34:D3:19:GLY:N	2.21	0.56
34:D3:54:LEU:O	34:D3:57:VAL:HG23	2.06	0.56
23:DB:1025:G:OP1	23:DB:1025:G:H8	1.89	0.56
23:DB:1447:C:H2'	23:DB:1448:G:H8	1.70	0.56
23:DB:1802:A:H2'	23:DB:1803:A:C8	2.41	0.56
23:DB:182:A:H1'	23:DB:434:U:H5'	1.86	0.56
23:DB:664:G:H2'	23:DB:665:U:H6	1.71	0.56
23:DB:692:C:H2'	23:DB:693:A:C8	2.41	0.56
23:DB:99:U:H5'	23:DB:100:U:OP1	2.05	0.56
25:DC:159:THR:O	25:DC:194:VAL:HG12	2.05	0.56
23:DB:691:C:H5'	25:DC:216:ARG:NH1	2.20	0.56
29:DE:146:VAL:O	29:DE:167:VAL:HA	2.06	0.56
47:DF:1:ALA:HB1	47:DF:4:HIS:HB3	1.87	0.56
47:DF:72:SER:HA	47:DF:78:ILE:HG22	1.87	0.56
40:DH:133:GLN:HG2	40:DH:139:PHE:HA	1.87	0.56
24:DI:1:ALA:CB	24:DI:2:LYS:HD2	2.35	0.56
41:DJ:25:LEU:HD22	41:DJ:26:GLY:H	1.70	0.56
43:DO:11:ALA:HB2	43:DO:96:GLY:CA	2.35	0.56
30:DY:15:ARG:N	30:DY:15:ARG:HD2	2.20	0.56
51:DZ:39:TRP:HB2	51:DZ:46:PHE:CE2	2.40	0.56
51:DZ:70:GLU:O	51:DZ:72:ARG:N	2.38	0.56
1:AA:106:C:H2'	1:AA:107:G:O4'	2.05	0.56
1:AA:1100:C:O2'	1:AA:1101:A:H5'	2.05	0.56
4:AE:131:ASN:ND2	4:AE:134:ASN:H	2.03	0.56
8:AI:103:VAL:HG23	8:AI:104:THR:N	2.20	0.56
8:AI:17:ARG:HB2	8:AI:65:THR:HB	1.87	0.56
8:AI:46:VAL:HA	8:AI:49:GLN:HG3	1.88	0.56
9:AJ:73:LEU:CD1	9:AJ:75:ASP:HB2	2.36	0.56
10:AK:62:ALA:HB1	10:AK:95:THR:HB	1.88	0.56
33:B1:34:GLU:CG	33:B1:49:LYS:HG3	2.35	0.56
23:BB:1932:A:H2'	23:BB:1933:G:O4'	2.05	0.56
23:BB:2537:U:H2'	23:BB:2538:C:H6	1.69	0.56
23:BB:2852:G:H2'	23:BB:2853:C:C6	2.41	0.56
23:BB:302:C:H2'	23:BB:303:G:H8	1.71	0.56
25:BC:20:ASN:HD21	25:BC:22:GLU:HG3	1.71	0.56
23:BB:2073:C:H5''	25:BC:227:VAL:HG12	1.87	0.56
25:BC:245:THR:C	25:BC:247:TRP:H	2.09	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:21:SER:HB3	27:BK:72:PRO:O	2.06	0.56
47:BF:108:PRO:C	47:BF:110:ILE:H	2.09	0.56
22:BA:43:C:H4'	47:BF:91:ARG:HG3	1.88	0.56
48:BG:70:LEU:O	48:BG:73:SER:HB3	2.05	0.56
40:BH:78:VAL:HG21	40:BH:142:VAL:HG13	1.88	0.56
43:BO:89:ASP:H	43:BO:115:LEU:HD13	1.69	0.56
28:BP:89:GLY:HA2	28:BP:111:GLU:C	2.26	0.56
28:BP:3:ILE:O	28:BP:7:LEU:HD13	2.06	0.56
1:CA:1041:G:H2'	1:CA:1042:A:H8	1.70	0.56
1:CA:124:C:O2'	1:CA:125:U:H5'	2.06	0.56
1:CA:1402:C:H2'	1:CA:1403:C:O4'	2.06	0.56
1:CA:462:G:H3'	1:CA:463:U:C6	2.40	0.56
1:CA:55:A:H2'	1:CA:56:U:O4'	2.06	0.56
1:CA:676:A:O2'	1:CA:677:U:H5'	2.05	0.56
1:CA:72:A:H2'	1:CA:73:C:O4'	2.05	0.56
1:CA:620:C:O2	3:CD:131:ILE:HG21	2.06	0.56
3:CD:9:LYS:HA	3:CD:12:ARG:HG3	1.86	0.56
12:CM:95:PRO:CA	12:CM:108:ARG:HG2	2.35	0.56
17:CT:81:GLN:O	17:CT:85:LEU:HD22	2.05	0.56
33:D1:34:GLU:CG	33:D1:49:LYS:HG3	2.34	0.56
22:DA:59:A:H2'	22:DA:60:C:C6	2.40	0.56
23:DB:1442:U:H2'	23:DB:1443:U:C6	2.41	0.56
23:DB:1851:U:H2'	23:DB:1852:U:H6	1.70	0.56
23:DB:1932:A:H2'	23:DB:1933:G:O4'	2.05	0.56
23:DB:2107:G:N3	23:DB:2107:G:H2'	2.19	0.56
23:DB:2157:G:N3	23:DB:2157:G:H2'	2.20	0.56
23:DB:41:C:H2'	23:DB:42:A:O4'	2.06	0.56
26:DD:101:PHE:O	26:DD:180:VAL:HG11	2.06	0.56
38:DM:34:LYS:HA	38:DM:100:LYS:O	2.06	0.56
45:DS:5:ALA:HB3	45:DS:54:ALA:HB2	1.87	0.56
45:DS:96:ILE:HG23	45:DS:96:ILE:O	2.05	0.56
30:DY:47:ILE:HD12	30:DY:54:VAL:HG21	1.87	0.56
1:AA:1002:G:H8	1:AA:1002:G:OP2	1.89	0.56
1:AA:1126:U:O2'	1:AA:1280:A:H2'	2.06	0.56
1:AA:1402:C:H2'	1:AA:1403:C:O4'	2.06	0.56
18:AB:125:PHE:N	18:AB:125:PHE:CD2	2.73	0.56
6:AG:145:GLU:HA	6:AG:148:LYS:HE3	1.88	0.56
15:AR:33:THR:HG23	15:AR:37:LYS:O	2.05	0.56
23:BB:1060:U:O4	24:BI:131:THR:HG22	2.05	0.56
23:BB:2140:G:N3	23:BB:2140:G:H2'	2.20	0.56
23:BB:2405:G:H1'	23:BB:2412:A:H61	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2542:A:H4'	23:BB:2543:G:H5'	1.88	0.56
23:BB:2662:A:H2'	23:BB:2663:G:O4'	2.06	0.56
23:BB:2742:G:O2'	23:BB:2743:U:H5'	2.06	0.56
27:BK:13:ASN:HD21	27:BK:98:ARG:HB2	1.69	0.56
42:BN:24:MET:SD	42:BN:44:LEU:HD22	2.45	0.56
42:BN:86:ARG:HB3	42:BN:117:ASP:OD1	2.06	0.56
43:BO:111:ARG:HG2	43:BO:117:PHE:H	1.70	0.56
43:BO:64:TYR:C	43:BO:66:GLY:H	2.09	0.56
28:BP:21:PRO:O	28:BP:46:VAL:HG23	2.06	0.56
44:BQ:91:ARG:HD3	49:BR:11:GLN:CD	2.26	0.56
45:BS:29:VAL:O	45:BS:33:LEU:HG	2.05	0.56
1:CA:1349:A:H2'	1:CA:1350:A:O4'	2.06	0.56
18:CB:86:CYS:H	18:CB:88:GLN:HE21	1.54	0.56
2:CC:154:GLY:HA3	2:CC:162:ALA:HB1	1.87	0.56
3:CD:36:ALA:HB2	3:CD:42:ALA:HB2	1.87	0.56
4:CE:158:LYS:NZ	7:CH:63:LYS:HD3	2.21	0.56
23:DB:1013:C:H2'	23:DB:1014:A:H8	1.70	0.56
23:DB:1911:U:H2'	23:DB:1918:A:N1	2.21	0.56
23:DB:2073:C:H5''	25:DC:227:VAL:HG12	1.88	0.56
23:DB:2243:U:H2'	23:DB:2244:U:C6	2.41	0.56
23:DB:2751:G:H3'	23:DB:2752:C:H6	1.71	0.56
23:DB:594:U:H2'	23:DB:595:C:H6	1.70	0.56
23:DB:633:A:H2'	23:DB:634:C:O4'	2.06	0.56
23:DB:1818:U:N3	25:DC:152:GLN:HB3	2.21	0.56
29:DE:146:VAL:HG11	29:DE:187:VAL:HG23	1.88	0.56
40:DH:77:THR:HA	40:DH:143:ILE:O	2.06	0.56
40:DH:3:VAL:HA	40:DH:37:VAL:O	2.06	0.56
24:DI:45:THR:CA	24:DI:48:ILE:HG22	2.35	0.56
41:DJ:81:ILE:HG12	41:DJ:82:GLY:N	2.19	0.56
28:DP:77:SER:OG	28:DP:79:VAL:HG22	2.06	0.56
44:DQ:20:ALA:HB2	44:DQ:38:VAL:CG2	2.36	0.56
46:DU:81:ARG:HD3	46:DU:96:LYS:HD2	1.87	0.56
35:DV:63:ILE:HD11	35:DV:72:VAL:HG22	1.87	0.56
30:DY:2:LYS:HD3	30:DY:2:LYS:H	1.70	0.56
1:AA:1071:C:O2'	1:AA:1072:G:H5'	2.05	0.56
1:AA:1110:A:H2'	1:AA:1111:A:H5'	1.88	0.56
1:AA:236:A:H2'	1:AA:237:G:H8	1.71	0.56
1:AA:704:A:H2'	1:AA:705:G:O4'	2.06	0.56
1:AA:733:G:O2'	1:AA:734:G:H5'	2.05	0.56
18:AB:187:ASP:OD1	18:AB:203:ASP:HB3	2.04	0.56
3:AD:52:VAL:HG12	3:AD:198:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:28:ASP:HB2	3:AD:33:ILE:HG21	1.88	0.56
5:AF:35:LYS:O	5:AF:64:VAL:HG13	2.06	0.56
6:AG:100:MET:O	6:AG:104:VAL:HG23	2.06	0.56
6:AG:73:GLU:O	6:AG:87:PRO:HA	2.06	0.56
9:AJ:42:LEU:HB3	9:AJ:71:LEU:HD21	1.88	0.56
11:AL:24:GLU:HB2	11:AL:26:CYS:HG	1.69	0.56
9:AJ:53:ILE:HG13	21:AN:84:ARG:CZ	2.36	0.56
16:AS:29:PRO:CB	16:AS:47:THR:HB	2.35	0.56
17:AT:61:ALA:HA	17:AT:66:ILE:HB	1.87	0.56
17:AT:81:GLN:O	17:AT:85:LEU:HD22	2.06	0.56
34:B3:14:LYS:HD3	34:B3:21:PHE:O	2.05	0.56
22:BA:87:U:H2'	22:BA:88:C:C5'	2.36	0.56
23:BB:1149:G:H2'	23:BB:1150:C:H6	1.69	0.56
23:BB:1442:U:H2'	23:BB:1443:U:C6	2.41	0.56
23:BB:3:U:O2'	23:BB:4:U:C6	2.59	0.56
23:BB:560:C:H2'	23:BB:561:G:O4'	2.06	0.56
23:BB:719:C:O2'	23:BB:720:U:H5'	2.06	0.56
25:BC:255:LYS:C	25:BC:257:ARG:H	2.10	0.56
23:BB:2578:G:N7	26:BD:145:SER:HB2	2.21	0.56
24:BI:76:ALA:O	24:BI:80:LYS:HG3	2.06	0.56
41:BJ:38:GLY:N	41:BJ:51:GLY:HA2	2.21	0.56
37:BL:57:LEU:C	37:BL:59:ARG:H	2.09	0.56
50:BT:32:LEU:H	50:BT:83:ALA:HB3	1.71	0.56
1:CA:1006:G:H2'	1:CA:1007:U:C6	2.40	0.56
1:CA:251:G:N3	1:CA:266:G:O6	2.39	0.56
11:CL:51:VAL:HG12	11:CL:52:CYS:N	2.19	0.56
21:CN:52:ARG:O	21:CN:55:SER:HB2	2.06	0.56
23:DB:1442:U:H2'	23:DB:1443:U:H6	1.70	0.56
23:DB:1854:A:H2	23:DB:2087:G:N3	2.04	0.56
23:DB:2359:C:H2'	23:DB:2360:G:C8	2.41	0.56
23:DB:309:A:N3	23:DB:329:G:O2'	2.39	0.56
23:DB:828:U:H2'	23:DB:829:A:C8	2.41	0.56
23:DB:828:U:H4'	23:DB:831:G:N1	2.20	0.56
47:DF:95:MET:O	47:DF:99:PHE:HB2	2.05	0.56
48:DG:70:LEU:O	48:DG:73:SER:HB3	2.06	0.56
24:DI:17:ALA:O	24:DI:18:ASN:CB	2.53	0.56
41:DJ:25:LEU:HA	41:DJ:28:LEU:HD22	1.87	0.56
42:DN:86:ARG:HB3	42:DN:117:ASP:OD1	2.06	0.56
42:DN:12:ARG:HG3	42:DN:13:ASN:H	1.70	0.56
43:DO:64:TYR:C	43:DO:66:GLY:H	2.10	0.56
43:DO:67:ASN:HB3	43:DO:70:ALA:CB	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DQ:84:LYS:HA	44:DQ:84:LYS:HE2	1.86	0.56
49:DR:37:GLU:O	49:DR:39:LEU:HD23	2.06	0.56
46:DU:12:VAL:HG22	46:DU:69:VAL:HG12	1.88	0.56
52:DW:48:ALA:O	52:DW:61:LYS:HB2	2.06	0.56
1:AA:1366:C:H2'	1:AA:1367:C:C6	2.41	0.56
1:AA:377:G:H2'	1:AA:378:G:H8	1.69	0.56
1:AA:772:U:H2'	1:AA:773:G:O4'	2.06	0.56
1:AA:844:G:N2	1:AA:845:A:H62	2.04	0.56
18:AB:107:ARG:NH2	18:AB:111:LYS:HB2	2.21	0.56
3:AD:117:VAL:HA	3:AD:122:ILE:CG1	2.35	0.56
6:AG:129:ASN:HA	6:AG:134:VAL:HG11	1.87	0.56
8:AI:11:ARG:HG3	8:AI:11:ARG:O	2.06	0.56
23:BB:1025:G:H8	23:BB:1025:G:OP1	1.88	0.56
23:BB:1150:C:H2'	23:BB:1151:A:H8	1.71	0.56
23:BB:1301:A:O2'	23:BB:1302:A:H3'	2.06	0.56
23:BB:1313:U:O2	23:BB:1313:U:H2'	2.06	0.56
23:BB:2134:A:N6	23:BB:2157:G:N2	2.54	0.56
23:BB:2230:G:H2'	23:BB:2231:U:H6	1.71	0.56
23:BB:584:C:OP1	44:BQ:5:ARG:HB3	2.06	0.56
23:BB:633:A:H2'	23:BB:634:C:O4'	2.05	0.56
23:BB:842:U:H2'	23:BB:843:G:H8	1.70	0.56
23:BB:969:G:OP1	30:BY:17:PRO:HG3	2.05	0.56
26:BD:3:GLY:C	26:BD:4:LEU:HD22	2.27	0.56
26:BD:31:ALA:HA	26:BD:97:SER:HA	1.87	0.56
29:BE:21:ARG:HG3	29:BE:22:ASP:O	2.06	0.56
41:BJ:81:ILE:HG12	41:BJ:82:GLY:N	2.20	0.56
37:BL:110:VAL:HG23	37:BL:126:ARG:O	2.05	0.56
37:BL:61:LEU:N	37:BL:61:LEU:HD12	2.21	0.56
35:BV:53:LYS:HA	35:BV:53:LYS:HZ3	1.71	0.56
18:CB:61:SER:HA	18:CB:224:ARG:HA	1.88	0.56
2:CC:18:ASN:ND2	2:CC:53:ARG:HH21	2.04	0.56
4:CE:101:GLY:H	4:CE:121:ASN:ND2	2.03	0.56
4:CE:14:LEU:HD22	4:CE:15:ILE:N	2.21	0.56
8:CI:32:ARG:HB3	8:CI:36:GLN:HE21	1.71	0.56
10:CK:126:ARG:HB2	19:CU:33:ARG:HD2	1.88	0.56
11:CL:43:LYS:HE3	11:CL:44:PRO:HD3	1.88	0.56
12:CM:38:ILE:HG22	12:CM:42:VAL:HG21	1.88	0.56
22:DA:95:U:H2'	22:DA:96:G:C8	2.40	0.56
23:DB:1505:A:H2'	23:DB:1506:U:C6	2.41	0.56
23:DB:2869:G:H2'	23:DB:2870:C:C6	2.41	0.56
23:DB:806:C:O2'	23:DB:807:U:H5'	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:55:LYS:HE2	26:DD:59:ARG:HB2	1.87	0.56
29:DE:116:ASP:O	29:DE:185:LYS:HE3	2.05	0.56
37:DL:80:SER:H	37:DL:113:ALA:CB	2.19	0.56
28:DP:6:GLN:O	28:DP:10:GLU:HB2	2.06	0.56
1:AA:1072:G:N2	18:AB:105:THR:HG21	2.21	0.55
1:AA:285:C:H2'	1:AA:286:C:C6	2.41	0.55
4:AE:155:LYS:HD2	7:AH:70:VAL:O	2.05	0.55
10:AK:80:ASN:CB	10:AK:105:ARG:HB3	2.35	0.55
1:AA:1227:A:H1'	12:AM:113:LYS:NZ	2.20	0.55
12:AM:79:LEU:HB2	12:AM:84:CYS:SG	2.47	0.55
13:AP:5:ARG:HH22	13:AP:24:SER:HA	1.72	0.55
23:BB:131:A:H2'	23:BB:132:G:C8	2.40	0.55
23:BB:1873:G:O2'	23:BB:1874:C:H5'	2.07	0.55
23:BB:1994:C:O2'	23:BB:1995:U:H5'	2.06	0.55
23:BB:2257:U:O2'	23:BB:2258:C:H5'	2.06	0.55
23:BB:664:G:H2'	23:BB:665:U:H6	1.71	0.55
26:BD:40:LEU:HA	26:BD:45:TYR:H	1.70	0.55
47:BF:134:GLN:C	47:BF:136:ILE:H	2.08	0.55
48:BG:83:THR:HA	48:BG:84:LYS:NZ	2.21	0.55
24:BI:105:LEU:HD11	24:BI:139:VAL:CG1	2.36	0.55
41:BJ:25:LEU:HA	41:BJ:28:LEU:HD22	1.87	0.55
27:BK:103:VAL:HG23	27:BK:122:VAL:O	2.06	0.55
43:BO:7:ARG:HA	43:BO:10:ARG:NE	2.21	0.55
28:BP:6:GLN:O	28:BP:10:GLU:HB2	2.06	0.55
46:BU:46:LYS:HZ2	46:BU:47:PRO:HD2	1.71	0.55
52:BW:43:LYS:HD2	52:BW:79:ILE:CD1	2.30	0.55
30:BY:6:ILE:HG22	30:BY:56:VAL:HG22	1.88	0.55
1:CA:1019:A:H2'	1:CA:1020:G:H8	1.71	0.55
1:CA:1408:A:O2'	1:CA:1409:C:H5'	2.06	0.55
1:CA:1507:A:H2'	1:CA:1508:A:C8	2.41	0.55
1:CA:1513:A:H2'	1:CA:1514:G:H8	1.69	0.55
1:CA:501:C:H2'	1:CA:502:A:C8	2.40	0.55
1:CA:663:A:O2'	1:CA:664:G:H5'	2.06	0.55
1:CA:908:A:O2'	1:CA:909:A:H5'	2.05	0.55
18:CB:18:GLN:HB2	18:CB:188:THR:OG1	2.07	0.55
2:CC:23:ALA:HB1	2:CC:27:GLU:OE1	2.06	0.55
5:CF:79:ARG:NH2	5:CF:87:SER:HB3	2.21	0.55
8:CI:90:ASP:C	8:CI:92:SER:H	2.09	0.55
1:CA:1060:U:OP1	9:CJ:53:ILE:HD11	2.06	0.55
12:CM:93:GLY:HA2	12:CM:108:ARG:HH12	1.71	0.55
23:DB:1081:U:H5'	24:DI:126:ARG:NH1	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:137:U:H3'	23:DB:138:U:C6	2.41	0.55
23:DB:1400:U:H2'	23:DB:1401:G:H8	1.71	0.55
23:DB:154:U:H2'	23:DB:155:A:H8	1.70	0.55
23:DB:1869:G:H3'	23:DB:1870:C:C5'	2.37	0.55
23:DB:2393:U:H2'	23:DB:2394:C:O4'	2.05	0.55
23:DB:370:G:O2'	23:DB:423:A:H3'	2.06	0.55
23:DB:547:A:H3'	23:DB:548:G:O4'	2.06	0.55
23:DB:553:G:O2'	23:DB:554:U:H5'	2.06	0.55
48:DG:84:LYS:C	48:DG:85:LYS:HG2	2.25	0.55
24:DI:5:GLN:O	24:DI:6:ALA:HB3	2.05	0.55
41:DJ:38:GLY:N	41:DJ:51:GLY:HA2	2.20	0.55
44:DQ:100:PHE:HB2	49:DR:13:ARG:HH12	1.71	0.55
45:DS:29:VAL:O	45:DS:33:LEU:HG	2.06	0.55
1:AA:1356:G:H2'	1:AA:1357:A:C8	2.42	0.55
1:AA:1507:A:H2'	1:AA:1508:A:C8	2.40	0.55
1:AA:624:C:H2'	1:AA:625:U:C6	2.41	0.55
1:AA:691:G:H1'	1:AA:696:A:N6	2.20	0.55
1:AA:926:G:H2'	1:AA:1505:G:N3	2.22	0.55
18:AB:85:SER:O	18:AB:86:CYS:HB2	2.06	0.55
11:AL:66:ILE:HG21	11:AL:71:HIS:HB3	1.88	0.55
12:AM:74:MET:O	12:AM:78:ARG:HB2	2.06	0.55
23:BB:1353:A:H2'	23:BB:1354:A:H8	1.71	0.55
23:BB:1430:G:H2'	23:BB:1431:A:C8	2.42	0.55
23:BB:1505:A:H2'	23:BB:1506:U:C6	2.41	0.55
23:BB:2405:G:H1'	23:BB:2412:A:N6	2.21	0.55
23:BB:2817:U:O2'	23:BB:2837:A:H1'	2.07	0.55
23:BB:674:G:H1'	29:BE:69:ARG:NE	2.09	0.55
23:BB:799:G:H3'	23:BB:800:A:H2'	1.87	0.55
25:BC:71:ASP:CB	25:BC:118:GLY:HA2	2.35	0.55
26:BD:124:ARG:HD3	26:BD:163:GLY:O	2.05	0.55
29:BE:149:ILE:O	29:BE:188:MET:HA	2.05	0.55
24:BI:91:LYS:HB2	24:BI:94:LYS:HD2	1.86	0.55
37:BL:109:LYS:HB2	37:BL:111:ILE:CD1	2.36	0.55
28:BP:4:ILE:HG22	28:BP:5:LYS:N	2.18	0.55
49:BR:5:PHE:HD1	49:BR:5:PHE:H	1.54	0.55
45:BS:24:ILE:HD11	45:BS:36:LEU:HD21	1.89	0.55
39:BX:21:LEU:HA	39:BX:25:GLN:HB3	1.88	0.55
51:BZ:70:GLU:O	51:BZ:72:ARG:N	2.39	0.55
1:CA:1242:G:O2'	1:CA:1243:C:H5'	2.06	0.55
1:CA:461:A:H3'	1:CA:462:G:O4'	2.06	0.55
1:CA:254:G:OP1	14:CQ:68:LYS:O	2.25	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CS:9:PHE:C	16:CS:10:ILE:HG13	2.27	0.55
23:DB:1051:G:H2'	23:DB:1052:C:C6	2.40	0.55
23:DB:1258:U:H2'	23:DB:1259:G:H8	1.71	0.55
23:DB:350:G:H2'	23:DB:351:C:O4'	2.06	0.55
23:DB:39:G:H2'	23:DB:40:U:H6	1.71	0.55
23:DB:720:U:H2'	23:DB:721:A:C8	2.42	0.55
23:DB:817:C:H2'	23:DB:818:G:O4'	2.06	0.55
23:DB:987:C:H2'	23:DB:988:A:O4'	2.06	0.55
40:DH:121:VAL:HG12	40:DH:123:ARG:HH11	1.71	0.55
23:DB:1252:G:H1'	44:DQ:32:ARG:HH22	1.70	0.55
49:DR:66:HIS:ND1	49:DR:94:THR:HG22	2.22	0.55
1:AA:1290:G:H2'	1:AA:1291:U:C6	2.41	0.55
1:AA:1329:A:OP1	12:AM:28:ARG:HB2	2.06	0.55
1:AA:251:G:N3	1:AA:266:G:O6	2.40	0.55
1:AA:97:G:H2'	1:AA:98:A:O4'	2.05	0.55
2:AC:168:ARG:HE	2:AC:169:GLU:N	2.04	0.55
3:AD:192:ALA:HB3	3:AD:194:ILE:HG22	1.87	0.55
5:AF:70:VAL:HG23	5:AF:71:ILE:N	2.22	0.55
6:AG:149:ALA:N	10:AK:55:ARG:NH2	2.54	0.55
8:AI:56:MET:HG3	8:AI:57:VAL:HG23	1.87	0.55
9:AJ:92:LEU:HD13	9:AJ:92:LEU:N	2.20	0.55
20:AO:11:ILE:CG1	20:AO:30:ALA:HB1	2.37	0.55
23:BB:1252:G:N2	44:BQ:32:ARG:HB3	2.22	0.55
23:BB:208:C:H2'	23:BB:209:C:C6	2.40	0.55
23:BB:2498:C:O2'	23:BB:2499:C:H5'	2.07	0.55
23:BB:2646:C:H2'	23:BB:2647:U:O4'	2.07	0.55
23:BB:348:A:H2'	23:BB:349:U:O4'	2.06	0.55
23:BB:737:C:O2'	23:BB:738:G:H5'	2.07	0.55
23:BB:936:A:H2'	23:BB:937:C:C6	2.41	0.55
25:BC:158:GLY:N	25:BC:194:VAL:HG13	2.21	0.55
25:BC:221:GLY:O	25:BC:223:ALA:N	2.33	0.55
23:BB:784:G:N1	25:BC:227:VAL:HG11	2.21	0.55
26:BD:102:ALA:H	26:BD:104:VAL:HG22	1.70	0.55
26:BD:114:LYS:HD2	26:BD:116:LYS:NZ	2.22	0.55
40:BH:90:LEU:HD11	40:BH:146:VAL:HG11	1.87	0.55
37:BL:132:ARG:HA	37:BL:135:ILE:HG22	1.87	0.55
1:CA:1512:U:H2'	1:CA:1513:A:C8	2.41	0.55
1:CA:195:A:H2'	1:CA:196:A:C8	2.41	0.55
1:CA:215:C:H2'	1:CA:216:U:H6	1.71	0.55
1:CA:242:G:H2'	1:CA:243:A:H5''	1.89	0.55
1:CA:304:U:H2'	1:CA:305:G:C8	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:476:U:O2'	1:CA:477:C:H5'	2.07	0.55
1:CA:648:A:H2'	1:CA:649:A:H8	1.71	0.55
1:CA:781:A:H2'	1:CA:782:A:H5'	1.88	0.55
18:CB:128:LEU:HD21	18:CB:132:GLU:HB3	1.88	0.55
4:CE:89:THR:HG21	4:CE:134:ASN:HD21	1.71	0.55
6:CG:63:VAL:C	6:CG:65:LEU:H	2.09	0.55
7:CH:24:VAL:HG12	7:CH:60:LEU:HB3	1.88	0.55
8:CI:19:PHE:O	8:CI:62:LEU:HA	2.06	0.55
9:CJ:15:HIS:H	9:CJ:70:HIS:CE1	2.24	0.55
10:CK:60:PHE:O	10:CK:64:VAL:HG12	2.05	0.55
12:CM:29:SER:HA	12:CM:32:ILE:CG2	2.35	0.55
23:DB:1018:U:O2'	23:DB:1019:U:H5'	2.07	0.55
23:DB:1061:U:H4'	23:DB:1070:A:O3'	2.06	0.55
23:DB:1513:U:O2'	23:DB:1514:G:H5'	2.06	0.55
23:DB:1859:U:H2'	23:DB:1860:G:C8	2.42	0.55
23:DB:191:A:H2'	23:DB:192:C:H6	1.70	0.55
23:DB:2256:G:H2'	23:DB:2257:U:H6	1.72	0.55
23:DB:2297:A:N1	23:DB:2320:U:H4'	2.21	0.55
23:DB:2452:C:H2'	23:DB:2453:A:C8	2.42	0.55
23:DB:319:G:H2'	23:DB:320:A:O4'	2.07	0.55
23:DB:842:U:H2'	23:DB:843:G:H8	1.70	0.55
23:DB:847:U:O4'	23:DB:847:U:O2	2.24	0.55
23:DB:1790:C:O2'	25:DC:207:ALA:HB2	2.06	0.55
26:DD:102:ALA:H	26:DD:104:VAL:HG22	1.70	0.55
26:DD:51:THR:HG22	26:DD:52:THR:H	1.71	0.55
48:DG:153:PRO:CG	48:DG:162:ARG:HB3	2.35	0.55
40:DH:131:SER:HB2	40:DH:141:LYS:CA	2.29	0.55
41:DJ:18:VAL:O	41:DJ:56:VAL:HA	2.06	0.55
37:DL:61:LEU:HD12	37:DL:61:LEU:N	2.21	0.55
43:DO:89:ASP:H	43:DO:115:LEU:HD13	1.70	0.55
23:DB:30:G:OP1	44:DQ:4:LYS:HG2	2.07	0.55
50:DT:39:THR:HG23	50:DT:41:ALA:H	1.70	0.55
30:DY:6:ILE:HG22	30:DY:56:VAL:HG22	1.89	0.55
1:AA:1221:G:P	16:AS:36:ARG:HH22	2.28	0.55
1:AA:151:A:H5'	1:AA:152:A:OP2	2.06	0.55
1:AA:242:G:H2'	1:AA:243:A:H5''	1.87	0.55
1:AA:499:A:H4'	1:AA:500:G:H5'	1.87	0.55
1:AA:842:U:H2'	1:AA:843:U:O3'	2.07	0.55
3:AD:89:LEU:HD22	3:AD:199:ILE:HD11	1.87	0.55
4:AE:47:PHE:O	4:AE:66:ALA:HA	2.07	0.55
6:AG:14:ASP:CB	6:AG:19:SER:H	2.12	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AJ:8:ILE:HD13	9:AJ:75:ASP:CA	2.36	0.55
10:AK:33:ILE:CB	10:AK:73:VAL:HG11	2.36	0.55
20:AO:69:TYR:CZ	20:AO:73:LYS:HG3	2.41	0.55
23:BB:1728:C:H2'	23:BB:1730:C:O2	2.06	0.55
23:BB:2728:U:H2'	23:BB:2729:G:C8	2.42	0.55
25:BC:132:ARG:HH22	25:BC:169:ALA:HA	1.71	0.55
25:BC:191:LEU:CD2	25:BC:192:GLY:H	2.18	0.55
29:BE:48:THR:HG22	29:BE:86:ALA:HB3	1.87	0.55
47:BF:8:LYS:HA	47:BF:12:VAL:CG2	2.36	0.55
40:BH:117:LEU:HG	40:BH:130:VAL:HG13	1.89	0.55
24:BI:109:ALA:HB1	24:BI:124:MET:HG3	1.89	0.55
24:BI:17:ALA:O	24:BI:18:ASN:HB3	2.06	0.55
44:BQ:20:ALA:HB2	44:BQ:38:VAL:CG2	2.36	0.55
44:BQ:91:ARG:HB3	44:BQ:93:ILE:HG22	1.87	0.55
50:BT:14:PRO:HA	50:BT:32:LEU:HB2	1.88	0.55
50:BT:44:LYS:O	50:BT:48:GLN:HG2	2.06	0.55
1:CA:1114:C:H2'	1:CA:1115:U:C6	2.41	0.55
1:CA:1253:G:N1	1:CA:1285:A:N6	2.54	0.55
1:CA:1332:A:H2'	1:CA:1333:A:H8	1.72	0.55
1:CA:1492:A:H5''	1:CA:1493:A:C4	2.42	0.55
1:CA:616:G:H2'	1:CA:616:G:N3	2.20	0.55
1:CA:980:C:H2'	1:CA:981:U:O4'	2.06	0.55
18:CB:15:PHE:O	18:CB:40:ILE:HD12	2.07	0.55
2:CC:72:PRO:HG2	2:CC:73:GLY:H	1.70	0.55
3:CD:117:VAL:HA	3:CD:122:ILE:CG1	2.37	0.55
3:CD:151:GLN:HB3	3:CD:153:ARG:HG2	1.87	0.55
5:CF:71:ILE:O	5:CF:75:GLU:HG3	2.06	0.55
1:CA:1228:C:OP1	12:CM:113:LYS:HB2	2.06	0.55
21:CN:40:ARG:NH1	21:CN:40:ARG:HB2	2.17	0.55
21:CN:9:GLU:HB2	21:CN:62:ARG:NH2	2.21	0.55
23:DB:1584:U:H5''	23:DB:1585:C:C5	2.41	0.55
23:DB:615:U:O4	29:DE:39:ALA:HB2	2.05	0.55
25:DC:4:LYS:CD	25:DC:5:CYS:H	2.19	0.55
47:DF:128:SER:HB3	47:DF:154:THR:HA	1.88	0.55
40:DH:119:ASN:N	40:DH:119:ASN:ND2	2.54	0.55
24:DI:129:GLU:HB3	24:DI:133:ARG:HH12	1.70	0.55
41:DJ:25:LEU:HD13	41:DJ:26:GLY:N	2.20	0.55
23:DB:637:A:OP2	37:DL:128:THR:HG21	2.06	0.55
37:DL:30:THR:O	37:DL:33:ARG:HG2	2.06	0.55
29:DE:29:HIS:NE2	37:DL:8:PRO:HG3	2.21	0.55
45:DS:15:GLN:O	45:DS:19:LEU:HB2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:32:LEU:O	50:DT:83:ALA:HB2	2.05	0.55
1:AA:1065:U:H1'	1:AA:1066:C:OP2	2.06	0.55
1:AA:184:G:N1	1:AA:194:C:N4	2.54	0.55
1:AA:231:U:H2'	1:AA:232:G:H8	1.70	0.55
1:AA:560:A:H4'	1:AA:561:U:C5'	2.36	0.55
18:AB:23:ASN:ND2	18:AB:25:LYS:HB2	2.19	0.55
1:AA:437:U:H5''	3:AD:151:GLN:HE21	1.70	0.55
8:AI:83:THR:HA	8:AI:86:LEU:HD22	1.89	0.55
9:AJ:52:LEU:HB2	21:AN:80:ARG:NE	2.21	0.55
12:AM:56:ARG:HH11	12:AM:56:ARG:HG3	1.72	0.55
12:AM:94:LEU:HB3	12:AM:95:PRO:HD2	1.89	0.55
13:AP:28:ARG:HH11	13:AP:29:ASN:HD22	1.53	0.55
22:BA:54:G:O2'	22:BA:55:U:H5'	2.07	0.55
23:BB:1041:G:O2'	23:BB:1042:G:H5'	2.07	0.55
23:BB:1258:U:H2'	23:BB:1259:G:H8	1.71	0.55
23:BB:1727:C:H2'	23:BB:1728:C:C6	2.42	0.55
23:BB:1777:U:O2'	23:BB:1778:U:H5'	2.06	0.55
23:BB:2183:A:H2'	23:BB:2184:A:C8	2.41	0.55
26:BD:136:ASN:HD21	26:BD:139:SER:C	2.10	0.55
47:BF:15:LEU:HA	47:BF:18:GLU:HB2	1.88	0.55
22:BA:42:C:C5	47:BF:65:LEU:HD22	2.41	0.55
48:BG:30:GLY:O	48:BG:78:VAL:HG12	2.07	0.55
27:BK:16:ALA:HB3	27:BK:47:ILE:HG13	1.89	0.55
37:BL:124:GLY:H	37:BL:143:GLU:HG3	1.68	0.55
49:BR:20:VAL:HG12	49:BR:21:ARG:H	1.71	0.55
30:BY:15:ARG:HD2	30:BY:15:ARG:N	2.22	0.55
30:BY:50:VAL:HB	30:BY:53:MET:CB	2.36	0.55
1:CA:121:U:H4'	1:CA:122:G:N7	2.20	0.55
1:CA:1258:G:H2'	1:CA:1259:C:C6	2.41	0.55
1:CA:253:A:H2'	1:CA:254:G:C8	2.42	0.55
1:CA:389:A:H3'	1:CA:390:U:H6	1.72	0.55
1:CA:843:U:H6	1:CA:843:U:H5''	1.71	0.55
3:CD:60:VAL:HA	3:CD:63:ILE:HD12	1.89	0.55
3:CD:8:LEU:O	3:CD:12:ARG:HG2	2.05	0.55
5:CF:66:ALA:HB1	5:CF:70:VAL:HG21	1.89	0.55
5:CF:84:VAL:HG22	5:CF:85:ILE:N	2.21	0.55
7:CH:95:MET:HB3	7:CH:99:GLY:H	1.71	0.55
13:CP:40:ASN:HD21	13:CP:43:ALA:N	2.05	0.55
31:D0:21:LEU:HD12	45:DS:19:LEU:O	2.06	0.55
23:DB:1116:G:O2'	23:DB:1117:C:H5'	2.06	0.55
23:DB:1461:C:H2'	23:DB:1462:C:C6	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1687:G:O2'	23:DB:1688:U:H5'	2.07	0.55
23:DB:1751:U:H2'	23:DB:1752:C:C5	2.41	0.55
23:DB:2261:C:O2'	23:DB:2262:U:H5'	2.06	0.55
23:DB:2555:U:H2'	23:DB:2556:C:O4'	2.07	0.55
23:DB:2798:U:H5''	23:DB:2799:A:OP1	2.06	0.55
23:DB:2813:A:H2'	23:DB:2814:A:C8	2.42	0.55
23:DB:2828:G:H2'	23:DB:2829:A:H8	1.72	0.55
23:DB:2849:U:N3	23:DB:2867:G:C8	2.74	0.55
23:DB:392:U:O2'	23:DB:393:C:H5'	2.07	0.55
29:DE:73:ILE:HG12	29:DE:73:ILE:O	2.06	0.55
47:DF:120:SER:HB2	47:DF:129:MET:HG3	1.88	0.55
27:DK:71:ARG:HA	27:DK:71:ARG:NE	2.22	0.55
50:DT:32:LEU:H	50:DT:83:ALA:HB3	1.72	0.55
46:DU:86:PHE:CG	46:DU:87:GLU:N	2.74	0.55
35:DV:41:GLU:C	35:DV:42:LEU:HD23	2.27	0.55
35:DV:70:ILE:HD13	35:DV:70:ILE:N	2.21	0.55
1:AA:1116:U:O2'	1:AA:1117:A:H5'	2.07	0.55
1:AA:1320:C:N3	16:AS:35:ARG:HD3	2.22	0.55
1:AA:1343:G:H2'	1:AA:1344:C:C6	2.42	0.55
1:AA:1463:U:H2'	1:AA:1464:U:C6	2.41	0.55
1:AA:1477:U:H2'	1:AA:1478:U:C6	2.41	0.55
1:AA:162:A:H2'	1:AA:163:C:O4'	2.06	0.55
1:AA:22:G:H2'	1:AA:23:C:H6	1.71	0.55
1:AA:613:C:H2'	1:AA:614:C:C6	2.42	0.55
1:AA:765:G:O6	1:AA:811:C:H5	1.90	0.55
18:AB:49:PHE:O	18:AB:53:LEU:HD13	2.06	0.55
3:AD:9:LYS:HA	3:AD:12:ARG:HG3	1.87	0.55
3:AD:29:THR:HG22	3:AD:30:LYS:HD3	1.88	0.55
11:AL:35:ARG:NH2	11:AL:36:VAL:HG22	2.19	0.55
20:AO:32:LEU:HD12	20:AO:63:ARG:HB2	1.88	0.55
33:B1:49:LYS:HG2	33:B1:50:GLU:N	2.17	0.55
23:BB:1115:G:H2'	23:BB:1116:G:C8	2.42	0.55
23:BB:1754:A:H2'	23:BB:1755:A:C8	2.41	0.55
23:BB:319:G:H2'	23:BB:320:A:O4'	2.07	0.55
23:BB:851:C:H2'	23:BB:852:U:H6	1.71	0.55
23:BB:93:G:H2'	23:BB:94:A:C8	2.42	0.55
26:BD:108:ASP:OD2	26:BD:173:GLN:HA	2.07	0.55
48:BG:36:LEU:N	48:BG:36:LEU:HD22	2.22	0.55
40:BH:128:HIS:O	40:BH:143:ILE:HG22	2.06	0.55
24:BI:10:LEU:HD12	24:BI:10:LEU:O	2.06	0.55
27:BK:25:LEU:CD2	27:BK:40:LYS:HB2	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:98:ILE:O	49:BR:99:THR:HG23	2.06	0.55
50:BT:39:THR:CG2	50:BT:42:GLU:H	2.20	0.55
52:BW:23:LYS:HD2	52:BW:24:ARG:N	2.21	0.55
1:CA:1269:A:H2'	1:CA:1270:G:H5'	1.89	0.55
1:CA:1339:A:H2'	1:CA:1340:A:O4'	2.06	0.55
1:CA:1366:C:O2'	1:CA:1367:C:H5'	2.06	0.55
1:CA:1389:C:H2'	1:CA:1390:U:H6	1.71	0.55
1:CA:978:A:H5'	1:CA:1362:A:H62	1.70	0.55
3:CD:54:LEU:HD22	3:CD:54:LEU:O	2.07	0.55
3:CD:90:LEU:HD21	3:CD:196:GLU:HB3	1.88	0.55
4:CE:47:PHE:O	4:CE:66:ALA:HA	2.06	0.55
5:CF:70:VAL:HG23	5:CF:71:ILE:N	2.21	0.55
20:CO:78:TYR:OH	20:CO:88:ARG:HG2	2.07	0.55
19:CU:36:PHE:HA	19:CU:39:LYS:HD2	1.89	0.55
22:DA:89:U:H5'	22:DA:90:C:C6	2.42	0.55
23:DB:256:A:H2'	23:DB:257:C:C6	2.41	0.55
23:DB:2676:C:H2'	23:DB:2677:G:H8	1.72	0.55
23:DB:2761:A:H1'	48:DG:142:GLN:NE2	2.21	0.55
23:DB:851:C:H2'	23:DB:852:U:C6	2.42	0.55
23:DB:915:C:H3'	23:DB:916:G:H8	1.71	0.55
23:DB:2787:C:H1'	26:DD:63:PRO:HG3	1.88	0.55
29:DE:161:ALA:HA	29:DE:164:LEU:HD12	1.88	0.55
22:DA:42:C:C5	47:DF:65:LEU:HD22	2.42	0.55
47:DF:69:ALA:HB3	47:DF:80:GLN:O	2.07	0.55
27:DK:76:VAL:H	28:DP:72:VAL:HG23	1.72	0.55
37:DL:57:LEU:C	37:DL:59:ARG:H	2.09	0.55
43:DO:52:SER:O	43:DO:58:ILE:HD12	2.07	0.55
46:DU:14:THR:O	46:DU:18:LYS:HG2	2.05	0.55
35:DV:6:ALA:HB3	35:DV:65:VAL:HG12	1.88	0.55
1:AA:1121:U:H2'	1:AA:1122:U:O4'	2.06	0.55
1:AA:469:C:H2'	1:AA:470:C:C6	2.42	0.55
18:AB:124:THR:OG1	18:AB:127:LYS:HE2	2.07	0.55
18:AB:214:GLY:O	18:AB:217:ALA:HB3	2.07	0.55
2:AC:149:LYS:HA	2:AC:168:ARG:HG3	1.88	0.55
1:AA:620:C:O2	3:AD:131:ILE:HG21	2.07	0.55
5:AF:3:HIS:CD2	5:AF:65:GLU:HG3	2.41	0.55
8:AI:44:ARG:O	8:AI:47:VAL:HG22	2.07	0.55
14:AQ:59:GLU:HB3	14:AQ:75:VAL:HG23	1.88	0.55
14:AQ:60:ILE:HG22	14:AQ:74:LEU:HA	1.89	0.55
23:BB:1637:A:H2'	23:BB:1638:C:C6	2.41	0.55
23:BB:176:A:O2'	23:BB:177:G:H5'	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1997:C:H2'	23:BB:1998:A:H8	1.71	0.55
47:BF:134:GLN:O	47:BF:136:ILE:N	2.40	0.55
40:BH:84:ALA:HA	40:BH:90:LEU:HG	1.88	0.55
1:CA:1226:C:N4	12:CM:102:LYS:HB2	2.22	0.55
1:CA:270:A:H2'	1:CA:271:C:H6	1.71	0.55
1:CA:295:C:H2'	1:CA:296:U:C6	2.41	0.55
1:CA:499:A:H4'	1:CA:500:G:OP1	2.05	0.55
18:CB:116:LEU:HD22	18:CB:140:LEU:HD11	1.88	0.55
3:CD:48:SER:O	3:CD:52:VAL:HG23	2.07	0.55
6:CG:68:VAL:HG11	6:CG:103:ILE:CG1	2.37	0.55
1:CA:1343:G:O3'	8:CI:123:ARG:HD3	2.07	0.55
12:CM:1:ALA:O	12:CM:3:ILE:HG13	2.07	0.55
20:CO:71:LYS:HZ3	20:CO:72:ARG:N	2.05	0.55
13:CP:53:ASP:O	13:CP:57:ILE:HG12	2.07	0.55
23:DB:1229:C:H2'	23:DB:1230:A:H8	1.72	0.55
23:DB:154:U:H2'	23:DB:155:A:C8	2.42	0.55
23:DB:2393:U:O2'	23:DB:2394:C:H5'	2.06	0.55
23:DB:2567:G:H2'	23:DB:2568:U:C6	2.41	0.55
23:DB:2769:U:H2'	23:DB:2770:G:H8	1.70	0.55
23:DB:2828:G:O2'	23:DB:2829:A:H5'	2.06	0.55
23:DB:2884:U:C6	31:D0:49:ARG:HG2	2.41	0.55
23:DB:419:U:H2'	23:DB:420:C:C6	2.42	0.55
23:DB:431:U:O2'	23:DB:432:A:H5'	2.07	0.55
23:DB:899:A:C2	23:DB:900:A:H1'	2.41	0.55
25:DC:146:LYS:HB3	25:DC:147:PRO:CD	2.32	0.55
25:DC:20:ASN:HD22	25:DC:23:LEU:HD13	1.71	0.55
25:DC:71:ASP:CB	25:DC:118:GLY:HA2	2.35	0.55
26:DD:33:ARG:NH1	26:DD:51:THR:HG22	2.22	0.55
47:DF:2:LYS:H	47:DF:2:LYS:HE3	1.72	0.55
48:DG:36:LEU:N	48:DG:36:LEU:HD22	2.21	0.55
40:DH:68:ARG:O	40:DH:72:ILE:HG12	2.07	0.55
23:DB:1006:C:H1'	41:DJ:108:MET:SD	2.47	0.55
38:DM:71:LYS:HG2	38:DM:93:VAL:HG12	1.89	0.55
49:DR:60:LYS:H	49:DR:100:GLY:HA3	1.70	0.55
51:DZ:32:ASN:HB2	51:DZ:53:ALA:H	1.70	0.55
1:AA:991:U:H3	1:AA:1212:U:C4'	2.19	0.55
1:AA:501:C:O2'	1:AA:502:A:H5'	2.06	0.55
1:AA:729:A:H2'	1:AA:730:G:H8	1.71	0.55
18:AB:53:LEU:HA	18:AB:56:LEU:HB2	1.88	0.55
7:AH:45:ILE:C	7:AH:63:LYS:HE3	2.26	0.55
8:AI:26:LYS:O	8:AI:61:ASP:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AP:22:ALA:HA	13:AP:33:ILE:HG13	1.88	0.55
33:B1:3:GLY:O	33:B1:5:ARG:N	2.40	0.55
23:BB:1315:C:H2'	23:BB:1316:U:H6	1.70	0.55
23:BB:1751:U:H2'	23:BB:1752:C:C5	2.42	0.55
23:BB:1771:C:H2'	23:BB:1772:A:C8	2.42	0.55
23:BB:1925:C:H2'	23:BB:1926:U:H5''	1.88	0.55
23:BB:256:A:H2'	23:BB:257:C:C6	2.41	0.55
23:BB:45:G:C5'	23:BB:46:G:H5'	2.29	0.55
23:BB:710:U:H2'	23:BB:711:G:H8	1.72	0.55
26:BD:33:ARG:HH11	26:BD:76:GLY:HA3	1.71	0.55
47:BF:69:ALA:HB3	47:BF:80:GLN:O	2.06	0.55
47:BF:87:LYS:CG	47:BF:88:VAL:H	2.19	0.55
48:BG:25:ILE:O	48:BG:32:LEU:HA	2.06	0.55
48:BG:89:VAL:HG23	48:BG:160:GLY:O	2.06	0.55
40:BH:116:ARG:HH22	40:BH:133:GLN:CA	2.19	0.55
27:BK:11:ALA:HB3	27:BK:85:VAL:HG22	1.87	0.55
27:BK:40:LYS:HZ2	27:BK:59:LYS:HE3	1.72	0.55
43:BO:58:ILE:HG22	43:BO:62:LEU:HD21	1.87	0.55
46:BU:86:PHE:CD1	46:BU:90:LYS:HB2	2.42	0.55
46:BU:86:PHE:HB3	46:BU:90:LYS:O	2.07	0.55
35:BV:63:ILE:CD1	35:BV:72:VAL:HG22	2.36	0.55
1:CA:1167:A:H2'	1:CA:1169:A:H62	1.72	0.55
1:CA:1271:A:H5'	1:CA:1314:C:H5'	1.89	0.55
1:CA:1411:C:O2'	1:CA:1412:C:H5'	2.07	0.55
1:CA:1478:U:H2'	1:CA:1479:C:C6	2.41	0.55
1:CA:239:U:C5'	1:CA:239:U:H6	2.20	0.55
1:CA:437:U:H5''	3:CD:151:GLN:HE21	1.72	0.55
1:CA:575:G:H4'	1:CA:576:C:O5'	2.07	0.55
1:CA:892:A:H2'	1:CA:893:C:C6	2.42	0.55
18:CB:8:MET:O	18:CB:10:LYS:N	2.39	0.55
3:CD:90:LEU:H	3:CD:90:LEU:HD22	1.72	0.55
4:CE:131:ASN:ND2	4:CE:134:ASN:H	2.05	0.55
6:CG:139:ASP:HA	6:CG:142:ARG:NH1	2.21	0.55
6:CG:94:ARG:O	6:CG:98:LEU:HB2	2.07	0.55
8:CI:108:ARG:NH1	8:CI:108:ARG:HB3	2.21	0.55
8:CI:62:LEU:HD22	8:CI:62:LEU:N	2.21	0.55
12:CM:2:ARG:HA	12:CM:7:ASN:O	2.07	0.55
1:CA:332:G:P	17:CT:2:ASN:HB3	2.47	0.55
19:CU:24:LYS:HD2	19:CU:25:ALA:N	2.21	0.55
22:DA:40:U:H1'	22:DA:43:C:C5	2.42	0.55
23:DB:1825:U:H2'	23:DB:1826:G:H8	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2292:U:H2'	23:DB:2293:G:C8	2.41	0.55
23:DB:292:U:O2'	23:DB:293:U:H5'	2.07	0.55
23:DB:37:C:O2'	23:DB:38:A:H5'	2.06	0.55
23:DB:402:A:C2'	23:DB:403:U:H5'	2.37	0.55
23:DB:670:A:H4'	23:DB:671:C:C5'	2.36	0.55
23:DB:719:C:O2'	23:DB:720:U:H5'	2.06	0.55
25:DC:177:SER:O	25:DC:270:ARG:HG3	2.06	0.55
47:DF:15:LEU:HA	47:DF:18:GLU:HB2	1.89	0.55
40:DH:135:HIS:HB3	40:DH:138:VAL:CB	2.35	0.55
27:DK:115:ILE:HG23	27:DK:116:ILE:N	2.22	0.55
27:DK:40:LYS:NZ	27:DK:59:LYS:HE3	2.22	0.55
46:DU:46:LYS:HZ2	46:DU:47:PRO:HD2	1.72	0.55
39:DX:21:LEU:HA	39:DX:25:GLN:HB3	1.89	0.55
23:DB:397:U:OP1	51:DZ:31:PRO:HA	2.07	0.55
1:AA:1396:A:C3'	1:AA:1397:C:H5'	2.36	0.55
1:AA:1412:C:H2'	1:AA:1413:A:H8	1.70	0.55
1:AA:767:A:H2'	1:AA:768:A:H8	1.72	0.55
18:AB:113:LEU:HD11	18:AB:144:GLU:HA	1.89	0.55
18:AB:79:VAL:HA	18:AB:213:LEU:HD21	1.88	0.55
2:AC:148:ILE:HA	2:AC:200:TRP:O	2.07	0.55
3:AD:156:ALA:O	3:AD:159:GLU:HB3	2.06	0.55
3:AD:36:ALA:HB2	3:AD:42:ALA:HB2	1.89	0.55
32:B4:8:LYS:HG2	32:B4:9:LYS:N	2.22	0.55
23:BB:154:U:H2'	23:BB:155:A:C8	2.41	0.55
23:BB:280:U:H2'	23:BB:281:C:C6	2.41	0.55
23:BB:480:A:N3	23:BB:480:A:H2'	2.22	0.55
23:BB:801:G:H4'	23:BB:802:A:OP2	2.07	0.55
23:BB:828:U:H2'	23:BB:829:A:C8	2.42	0.55
23:BB:909:A:H2'	23:BB:912:C:H5	1.72	0.55
25:BC:146:LYS:HB3	25:BC:147:PRO:CD	2.31	0.55
40:BH:100:ALA:O	40:BH:110:VAL:HG23	2.07	0.55
40:BH:42:LYS:HA	40:BH:46:PHE:HB2	1.87	0.55
41:BJ:19:ASP:HB2	41:BJ:57:LEU:HB2	1.89	0.55
37:BL:30:THR:O	37:BL:33:ARG:HG2	2.05	0.55
42:BN:2:ARG:HB3	42:BN:2:ARG:CZ	2.37	0.55
42:BN:34:ILE:HG22	42:BN:35:LYS:N	2.22	0.55
28:BP:3:ILE:HG23	28:BP:4:ILE:H	1.72	0.55
50:BT:30:ILE:HG23	50:BT:85:VAL:CG2	2.37	0.55
46:BU:5:ARG:NH2	46:BU:93:ARG:HD3	2.21	0.55
52:BW:37:VAL:HB	52:BW:38:ARG:HD3	1.88	0.55
1:CA:1355:G:O2'	1:CA:1356:G:H5'	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CK:95:THR:HG23	10:CK:96:ILE:N	2.22	0.55
14:CQ:59:GLU:HB3	14:CQ:75:VAL:CG2	2.37	0.55
23:DB:1015:U:H2'	23:DB:1016:G:C8	2.42	0.55
23:DB:103:A:H3'	23:DB:104:A:H8	1.72	0.55
23:DB:1794:A:H2'	23:DB:1795:C:C6	2.41	0.55
23:DB:18:U:H2'	23:DB:19:A:C8	2.42	0.55
23:DB:394:C:O2'	23:DB:395:U:H5'	2.07	0.55
23:DB:551:G:H2'	23:DB:552:U:H6	1.72	0.55
47:DF:110:ILE:CG2	47:DF:113:PHE:HB3	2.37	0.55
47:DF:137:PHE:HB2	47:DF:138:PRO:CD	2.36	0.55
43:DO:76:LYS:O	43:DO:79:ALA:HB3	2.06	0.55
1:AA:203:G:H5'	1:AA:468:A:O2'	2.06	0.55
1:AA:55:A:H2'	1:AA:56:U:O4'	2.07	0.55
1:AA:749:A:O2'	1:AA:750:C:H5'	2.06	0.55
1:AA:802:A:H2'	1:AA:803:G:O4'	2.07	0.55
18:AB:112:ARG:HE	18:AB:116:LEU:CD1	2.20	0.55
18:AB:216:VAL:O	18:AB:220:VAL:HG23	2.07	0.55
18:AB:55:GLU:O	18:AB:59:ILE:HG12	2.07	0.55
3:AD:29:THR:HB	3:AD:30:LYS:HZ3	1.70	0.55
6:AG:147:ASN:HA	10:AK:55:ARG:CZ	2.37	0.55
33:B1:36:LYS:CB	33:B1:47:ILE:HA	2.37	0.55
23:BB:1309:G:H4'	36:B2:7:PRO:HB2	1.89	0.55
23:BB:1210:G:H5'	23:BB:1212:G:O4'	2.07	0.55
23:BB:1818:U:N3	25:BC:152:GLN:HB3	2.20	0.55
23:BB:1967:C:H2'	23:BB:1968:G:O4'	2.07	0.55
23:BB:68:G:O2'	23:BB:69:C:H5'	2.06	0.55
26:BD:178:VAL:HG12	26:BD:179:ARG:HG3	1.89	0.55
29:BE:148:ILE:HA	29:BE:187:VAL:HB	1.87	0.55
29:BE:2:GLU:HA	29:BE:13:THR:CA	2.37	0.55
23:BB:518:G:H4'	45:BS:18:ARG:NH2	2.22	0.55
23:BB:923:G:N3	52:BW:23:LYS:HE3	2.22	0.55
1:CA:1286:U:O2	1:CA:1286:U:O4'	2.25	0.55
1:CA:13:U:O2	1:CA:914:A:H3'	2.07	0.55
1:CA:151:A:H5'	1:CA:152:A:OP2	2.06	0.55
1:CA:201:G:H2'	1:CA:202:G:O4'	2.07	0.55
1:CA:203:G:H5'	1:CA:468:A:O2'	2.06	0.55
1:CA:364:A:H2'	1:CA:365:U:O2	2.07	0.55
18:CB:107:ARG:HA	18:CB:110:ILE:CD1	2.36	0.55
3:CD:32:LYS:HA	3:CD:35:GLN:HE21	1.72	0.55
6:CG:63:VAL:HA	6:CG:66:GLU:OE2	2.07	0.55
11:CL:49:ARG:HH12	11:CL:88:ASP:CB	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1329:A:H4'	12:CM:23:GLY:O	2.06	0.55
36:D2:21:ARG:HD3	36:D2:30:VAL:HG11	1.89	0.55
23:DB:2322:A:H3'	23:DB:2323:G:H8	1.72	0.55
23:DB:237:C:O2'	23:DB:238:C:H5'	2.06	0.55
23:DB:2455:G:H2'	23:DB:2456:C:C6	2.41	0.55
23:DB:2885:G:H2'	23:DB:2886:A:O4'	2.06	0.55
23:DB:45:G:C5'	23:DB:46:G:H5'	2.31	0.55
23:DB:854:C:O2'	23:DB:855:G:H5'	2.07	0.55
23:DB:863:A:H2'	23:DB:864:G:H8	1.71	0.55
25:DC:83:ASP:HB2	25:DC:90:ILE:HG12	1.88	0.55
47:DF:35:LEU:HA	47:DF:153:ILE:HG12	1.89	0.55
48:DG:8:VAL:CG1	48:DG:49:LEU:HB3	2.36	0.55
24:DI:71:LYS:HB3	24:DI:115:ASP:OD2	2.06	0.55
23:DB:2358:A:H61	37:DL:54:GLN:HE22	1.55	0.55
30:DY:6:ILE:N	30:DY:6:ILE:HD13	2.20	0.55
1:AA:1014:A:C5'	16:AS:13:HIS:HB3	2.37	0.54
1:AA:213:G:H3'	1:AA:214:C:H6	1.72	0.54
1:AA:499:A:H4'	1:AA:500:G:OP1	2.07	0.54
1:AA:630:A:H2'	1:AA:631:C:C6	2.42	0.54
5:AF:56:LYS:N	5:AF:56:LYS:HE3	2.22	0.54
6:AG:50:ALA:HB2	6:AG:57:GLU:HG3	1.89	0.54
8:AI:97:LEU:HA	8:AI:102:PHE:HD1	1.73	0.54
10:AK:35:ASP:OD2	10:AK:37:GLN:HG2	2.07	0.54
11:AL:3:VAL:O	11:AL:7:VAL:HG23	2.08	0.54
20:AO:85:LEU:N	20:AO:85:LEU:HD12	2.22	0.54
34:B3:7:ARG:HH11	34:B3:7:ARG:HG3	1.72	0.54
23:BB:1038:G:H2'	23:BB:1039:A:C8	2.42	0.54
23:BB:1400:U:H2'	23:BB:1401:G:H8	1.71	0.54
23:BB:1479:G:O2'	23:BB:1480:C:H5'	2.07	0.54
23:BB:2181:U:H2'	23:BB:2182:U:C6	2.42	0.54
23:BB:2730:C:H2'	23:BB:2731:G:H8	1.71	0.54
23:BB:441:U:H2'	23:BB:442:G:C8	2.42	0.54
23:BB:553:G:O2'	23:BB:554:U:H5'	2.06	0.54
23:BB:832:U:H2'	23:BB:833:A:H8	1.70	0.54
25:BC:155:ARG:CB	25:BC:155:ARG:HH11	2.19	0.54
29:BE:157:LEU:HG	29:BE:169:VAL:HG11	1.88	0.54
29:BE:34:ALA:CB	29:BE:96:VAL:HG21	2.35	0.54
40:BH:89:LYS:HZ2	40:BH:89:LYS:HA	1.71	0.54
24:BI:14:ALA:HB1	24:BI:50:LYS:HA	1.88	0.54
27:BK:85:VAL:HG21	27:BK:115:ILE:CD1	2.36	0.54
43:BO:76:LYS:O	43:BO:79:ALA:HB3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BU:64:ILE:HG13	46:BU:65:GLN:N	2.22	0.54
46:BU:12:VAL:HA	46:BU:69:VAL:HG12	1.87	0.54
1:CA:1122:U:H2'	1:CA:1123:U:C6	2.41	0.54
1:CA:1127:G:H2'	1:CA:1128:C:C6	2.41	0.54
1:CA:1340:A:H2'	1:CA:1341:U:O4'	2.07	0.54
1:CA:1389:C:H2'	1:CA:1390:U:C6	2.41	0.54
1:CA:202:G:H2'	1:CA:203:G:H8	1.72	0.54
1:CA:237:G:H5''	14:CQ:26:ARG:HH21	1.72	0.54
1:CA:376:G:H4'	13:CP:5:ARG:HD3	1.89	0.54
18:CB:75:ALA:O	18:CB:79:VAL:HG23	2.07	0.54
5:CF:3:HIS:CD2	5:CF:65:GLU:HG3	2.41	0.54
12:CM:78:ARG:NH1	16:CS:64:GLU:HG2	2.22	0.54
13:CP:6:LEU:CD1	13:CP:71:VAL:HB	2.37	0.54
22:DA:37:C:H2'	22:DA:38:C:O4'	2.06	0.54
23:DB:1144:A:O2'	23:DB:1145:C:H5'	2.07	0.54
23:DB:2231:U:H2'	23:DB:2232:C:H6	1.72	0.54
23:DB:2556:C:H2'	23:DB:2557:G:O4'	2.07	0.54
24:DI:2:LYS:NZ	24:DI:2:LYS:HB3	2.22	0.54
23:DB:1666:G:O3'	27:DK:6:THR:HG23	2.07	0.54
35:DV:63:ILE:CD1	35:DV:72:VAL:HG22	2.36	0.54
52:DW:17:ALA:O	52:DW:18:LYS:HD2	2.07	0.54
1:AA:1305:G:H1'	1:AA:1332:A:N6	2.21	0.54
1:AA:409:U:H2'	1:AA:410:G:C8	2.42	0.54
18:AB:138:ARG:HA	18:AB:141:GLU:OE1	2.06	0.54
18:AB:160:LEU:HB3	18:AB:182:VAL:HG22	1.88	0.54
18:AB:221:ARG:HG3	18:AB:222:GLU:N	2.22	0.54
2:AC:13:ILE:HG12	2:AC:14:VAL:HG13	1.89	0.54
4:AE:103:GLY:O	4:AE:121:ASN:HA	2.08	0.54
9:AJ:18:ILE:HG12	9:AJ:72:ARG:HG2	1.89	0.54
12:AM:80:MET:HE2	12:AM:90:HIS:HB2	1.90	0.54
1:AA:668:G:O2'	20:AO:46:HIS:HB3	2.07	0.54
23:BB:1042:G:O2'	23:BB:1043:C:H5'	2.07	0.54
23:BB:1400:U:H2'	23:BB:1401:G:C8	2.41	0.54
23:BB:1734:G:O2'	23:BB:1735:A:H5'	2.07	0.54
23:BB:182:A:H1'	23:BB:434:U:H5'	1.88	0.54
23:BB:2556:C:H2'	23:BB:2557:G:O4'	2.06	0.54
23:BB:431:U:O2'	23:BB:432:A:H5'	2.08	0.54
23:BB:594:U:H2'	23:BB:595:C:H6	1.73	0.54
23:BB:759:G:H2'	23:BB:760:G:C8	2.42	0.54
23:BB:955:U:H5''	38:BM:86:LYS:HZ2	1.72	0.54
26:BD:107:VAL:H	26:BD:205:PRO:HA	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BI:5:GLN:HG2	24:BI:6:ALA:N	2.22	0.54
27:BK:2:ILE:HG22	27:BK:32:TYR:HB3	1.90	0.54
42:BN:57:THR:HG22	42:BN:58:ASP:H	1.71	0.54
45:BS:42:LYS:O	45:BS:45:VAL:HG22	2.06	0.54
45:BS:96:ILE:HG23	45:BS:96:ILE:O	2.07	0.54
35:BV:63:ILE:HD12	35:BV:63:ILE:N	2.21	0.54
1:CA:1432:G:H1'	1:CA:1468:A:N6	2.22	0.54
18:CB:130:LYS:HA	18:CB:130:LYS:HE2	1.88	0.54
18:CB:26:MET:HG3	18:CB:188:THR:O	2.06	0.54
2:CC:18:ASN:CG	2:CC:53:ARG:HH21	2.10	0.54
2:CC:71:ARG:HB2	2:CC:74:ILE:CG2	2.38	0.54
12:CM:69:ARG:HH11	12:CM:69:ARG:HA	1.73	0.54
32:D4:8:LYS:HG2	32:D4:9:LYS:N	2.22	0.54
22:DA:55:U:H2'	22:DA:56:G:H8	1.71	0.54
23:DB:1244:A:O2'	23:DB:1245:G:H5'	2.07	0.54
23:DB:1558:C:H4'	23:DB:1559:U:C5'	2.35	0.54
23:DB:189:G:H2'	23:DB:205:G:H22	1.72	0.54
23:DB:2405:G:H1'	23:DB:2412:A:N6	2.23	0.54
23:DB:2438:U:H2'	23:DB:2441:U:H5	1.71	0.54
23:DB:417:C:H2'	23:DB:418:C:H6	1.72	0.54
23:DB:750:A:H2'	23:DB:751:A:H5''	1.89	0.54
26:DD:33:ARG:NE	26:DD:51:THR:HB	2.21	0.54
47:DF:79:ARG:H	47:DF:82:TYR:HB2	1.72	0.54
24:DI:49:GLU:CB	24:DI:52:LEU:HD12	2.38	0.54
42:DN:19:ALA:C	42:DN:21:PHE:H	2.10	0.54
23:DB:2816:G:O3'	42:DN:99:LYS:HE3	2.07	0.54
1:AA:1315:U:H5'	1:AA:1316:G:OP2	2.08	0.54
1:AA:332:G:P	17:AT:2:ASN:HB3	2.47	0.54
1:AA:335:C:H2'	1:AA:336:A:C8	2.42	0.54
1:AA:50:A:H1'	1:AA:52:C:C6	2.42	0.54
1:AA:865:A:H2	1:AA:918:A:H4'	1.71	0.54
1:AA:900:A:H2'	1:AA:901:A:H8	1.71	0.54
18:AB:83:ALA:HA	18:AB:88:GLN:NE2	2.22	0.54
2:AC:26:LYS:HG3	2:AC:27:GLU:N	2.15	0.54
8:AI:11:ARG:HG2	8:AI:77:ALA:CB	2.38	0.54
10:AK:60:PHE:O	10:AK:64:VAL:HG12	2.07	0.54
12:AM:105:ALA:O	12:AM:109:LYS:HG2	2.08	0.54
21:AN:50:LEU:N	21:AN:51:PRO:CD	2.66	0.54
14:AQ:68:LYS:HG2	14:AQ:69:THR:HG23	1.88	0.54
22:BA:89:U:H5'	22:BA:90:C:C6	2.42	0.54
23:BB:1506:U:H2'	23:BB:1507:C:C6	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1645:G:H5''	23:BB:1646:C:H5'	1.90	0.54
23:BB:39:G:H2'	23:BB:40:U:C6	2.42	0.54
23:BB:585:G:H2'	23:BB:1251:C:H42	1.71	0.54
23:BB:598:U:H2'	23:BB:599:A:H8	1.72	0.54
23:BB:639:U:H2'	23:BB:640:C:C6	2.42	0.54
23:BB:67:U:H2'	23:BB:68:G:C8	2.41	0.54
23:BB:847:U:O4'	23:BB:847:U:O2	2.24	0.54
29:BE:161:ALA:HA	29:BE:164:LEU:HD12	1.90	0.54
47:BF:137:PHE:HB2	47:BF:138:PRO:CD	2.36	0.54
48:BG:3:VAL:O	48:BG:68:ARG:HG3	2.08	0.54
26:BD:20:VAL:HA	27:BK:72:PRO:HB3	1.90	0.54
42:BN:22:ARG:CZ	42:BN:70:THR:H	2.20	0.54
43:BO:11:ALA:HB2	43:BO:96:GLY:CA	2.36	0.54
43:BO:76:LYS:O	43:BO:80:GLU:HG2	2.07	0.54
23:BB:141:G:H1	50:BT:1:MET:HA	1.72	0.54
46:BU:81:ARG:HD3	46:BU:96:LYS:HD2	1.87	0.54
39:BX:23:ARG:HD2	39:BX:27:ASN:ND2	2.22	0.54
1:CA:161:A:H2'	1:CA:162:A:C8	2.42	0.54
1:CA:202:G:H2'	1:CA:203:G:C8	2.42	0.54
1:CA:45:G:H2'	1:CA:46:G:H8	1.72	0.54
2:CC:63:ILE:HD12	2:CC:98:ALA:HB2	1.90	0.54
4:CE:95:MET:HA	4:CE:124:ALA:CB	2.38	0.54
10:CK:80:ASN:CB	10:CK:105:ARG:HB3	2.37	0.54
36:D2:3:ARG:NE	36:D2:3:ARG:HA	2.22	0.54
34:D3:14:LYS:O	34:D3:21:PHE:HA	2.08	0.54
23:DB:106:C:H2'	23:DB:107:G:H8	1.72	0.54
1:CA:1517:G:H1'	23:DB:1919:A:O3'	2.07	0.54
23:DB:315:G:H2'	23:DB:316:C:C6	2.43	0.54
23:DB:876:C:H42	23:DB:901:C:H41	1.53	0.54
26:DD:5:VAL:HG23	26:DD:32:ASN:HD21	1.72	0.54
47:DF:87:LYS:HG3	47:DF:88:VAL:H	1.72	0.54
48:DG:51:PHE:CE2	48:DG:68:ARG:HA	2.42	0.54
37:DL:3:LEU:O	37:DL:5:THR:N	2.39	0.54
38:DM:17:ASN:HD21	38:DM:95:LEU:CD1	2.20	0.54
52:DW:17:ALA:CA	52:DW:35:ILE:HG23	2.27	0.54
1:AA:1221:G:H5''	1:AA:1321:U:O2	2.07	0.54
1:AA:167:A:O2'	1:AA:168:G:H5'	2.07	0.54
18:AB:127:LYS:HB3	18:AB:127:LYS:NZ	2.22	0.54
2:AC:48:LYS:H	2:AC:48:LYS:HD3	1.72	0.54
6:AG:30:MET:HG3	6:AG:38:ALA:HB2	1.90	0.54
9:AJ:8:ILE:HB	9:AJ:73:LEU:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AM:72:ILE:HG22	12:AM:76:ILE:HD11	1.89	0.54
21:AN:59:GLN:HE21	21:AN:59:GLN:N	2.04	0.54
13:AP:6:LEU:CD1	13:AP:71:VAL:HB	2.38	0.54
22:BA:40:U:H1'	22:BA:43:C:C5	2.41	0.54
23:BB:1180:U:H2'	23:BB:1181:U:C6	2.43	0.54
23:BB:2333:A:H5'	23:BB:2335:A:H1'	1.89	0.54
23:BB:2784:U:H2'	23:BB:2785:C:C6	2.41	0.54
23:BB:607:U:O4	23:BB:620:G:H5''	2.08	0.54
23:BB:987:C:H2'	23:BB:988:A:O4'	2.08	0.54
25:BC:20:ASN:HD22	25:BC:23:LEU:HD13	1.73	0.54
48:BG:93:TYR:C	48:BG:94:ARG:HG3	2.27	0.54
27:BK:109:SER:C	27:BK:111:LYS:H	2.11	0.54
39:BX:39:GLN:O	39:BX:42:LEU:HB2	2.07	0.54
1:CA:1356:G:H2'	1:CA:1357:A:H8	1.72	0.54
1:CA:149:A:H1'	1:CA:1446:A:C2	2.42	0.54
1:CA:434:U:H3'	1:CA:435:A:H8	1.72	0.54
1:CA:880:C:H2'	1:CA:881:G:C8	2.41	0.54
2:CC:34:SER:O	2:CC:38:VAL:HG22	2.07	0.54
3:CD:28:ASP:HB2	3:CD:33:ILE:HG21	1.89	0.54
3:CD:78:ALA:HA	3:CD:81:LEU:HD12	1.90	0.54
4:CE:140:ILE:HG22	4:CE:144:GLU:OE1	2.07	0.54
4:CE:83:PRO:HB3	4:CE:96:GLN:HG3	1.90	0.54
6:CG:74:VAL:HA	6:CG:87:PRO:HA	1.89	0.54
7:CH:28:SER:HB2	7:CH:56:PRO:HB2	1.88	0.54
10:CK:35:ASP:OD2	10:CK:37:GLN:HG2	2.07	0.54
11:CL:66:ILE:HG21	11:CL:71:HIS:HB3	1.88	0.54
22:DA:50:A:OP1	43:DO:68:LYS:HB2	2.08	0.54
23:DB:585:G:H2'	23:DB:1251:C:H42	1.72	0.54
23:DB:1400:U:H2'	23:DB:1401:G:C8	2.41	0.54
23:DB:1414:C:H2'	23:DB:1415:U:C6	2.43	0.54
23:DB:1520:U:H2'	23:DB:1521:G:O4'	2.07	0.54
23:DB:19:A:H2'	23:DB:20:C:C6	2.43	0.54
23:DB:2095:A:H3'	23:DB:2096:C:H6	1.71	0.54
23:DB:2063:C:O2	23:DB:2450:A:N1	2.39	0.54
23:DB:2458:G:N3	23:DB:2458:G:H2'	2.22	0.54
23:DB:2599:G:H2'	23:DB:2600:A:H8	1.72	0.54
23:DB:2645:G:C5'	23:DB:2732:G:H8	2.20	0.54
23:DB:506:G:H5'	23:DB:509:C:H1'	1.89	0.54
23:DB:674:G:H5''	29:DE:71:GLY:N	2.22	0.54
23:DB:742:A:H2'	23:DB:743:A:C8	2.40	0.54
25:DC:28:PRO:HG2	25:DC:33:LEU:HD11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:109:ARG:HB3	47:DF:135:ILE:HB	1.90	0.54
40:DH:57:LYS:HG3	40:DH:58:LEU:HD23	1.90	0.54
27:DK:25:LEU:HD21	27:DK:40:LYS:N	2.21	0.54
27:DK:99:ILE:HB	27:DK:118:LEU:HD22	1.90	0.54
49:DR:20:VAL:HG12	49:DR:21:ARG:H	1.72	0.54
50:DT:57:VAL:HG13	50:DT:58:VAL:N	2.23	0.54
46:DU:94:PHE:HA	46:DU:101:THR:HA	1.89	0.54
46:DU:82:VAL:H	46:DU:96:LYS:NZ	2.04	0.54
1:AA:389:A:H3'	1:AA:390:U:H6	1.73	0.54
1:AA:489:C:H2'	1:AA:490:C:C6	2.43	0.54
1:AA:957:U:H2'	1:AA:959:A:OP2	2.07	0.54
3:AD:56:GLU:O	3:AD:60:VAL:HG12	2.06	0.54
4:AE:92:ARG:HB3	4:AE:92:ARG:NH1	2.23	0.54
7:AH:24:VAL:HG12	7:AH:60:LEU:HB3	1.88	0.54
9:AJ:57:VAL:HG22	9:AJ:58:ASN:N	2.18	0.54
1:AA:376:G:H4'	13:AP:5:ARG:HD3	1.89	0.54
23:BB:1723:G:H3'	23:BB:1724:G:C8	2.38	0.54
23:BB:172:A:H2'	23:BB:173:A:H8	1.72	0.54
23:BB:1802:A:H2'	23:BB:1803:A:C8	2.42	0.54
23:BB:2471:A:O2'	23:BB:2472:G:C8	2.59	0.54
23:BB:2798:U:H5'	23:BB:2799:A:OP1	2.06	0.54
23:BB:720:U:H2'	23:BB:721:A:C8	2.41	0.54
25:BC:78:GLU:HB2	25:BC:92:LEU:HG	1.89	0.54
23:BB:2579:C:O2'	26:BD:136:ASN:HA	2.06	0.54
47:BF:102:LEU:HD22	47:BF:106:ALA:CB	2.37	0.54
47:BF:33:ILE:HB	47:BF:90:LEU:HB2	1.88	0.54
47:BF:79:ARG:H	47:BF:82:TYR:HB2	1.72	0.54
40:BH:73:ASN:CB	40:BH:141:LYS:HB2	2.38	0.54
40:BH:73:ASN:CB	40:BH:141:LYS:HE2	2.37	0.54
41:BJ:25:LEU:HD13	41:BJ:26:GLY:N	2.21	0.54
27:BK:7:MET:SD	27:BK:20:MET:HB2	2.47	0.54
38:BM:42:THR:H	38:BM:45:GLN:CD	2.11	0.54
50:BT:39:THR:HG23	50:BT:41:ALA:H	1.72	0.54
46:BU:14:THR:O	46:BU:18:LYS:HG2	2.07	0.54
46:BU:94:PHE:HA	46:BU:101:THR:HA	1.88	0.54
52:BW:49:ASN:CB	52:BW:60:ALA:HA	2.37	0.54
1:CA:1040:U:H2'	1:CA:1041:G:O4'	2.07	0.54
1:CA:1174:G:O2'	1:CA:1175:G:H5'	2.08	0.54
1:CA:363:A:P	11:CL:57:THR:HG21	2.48	0.54
1:CA:484:G:H5'	1:CA:486:U:H5'	1.89	0.54
1:CA:539:A:H2'	1:CA:540:G:C8	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:729:A:H2'	1:CA:730:G:C8	2.41	0.54
1:CA:91:U:H2'	1:CA:92:U:C6	2.43	0.54
18:CB:65:LYS:O	18:CB:66:ILE:HD13	2.08	0.54
18:CB:93:HIS:O	18:CB:94:ARG:HG2	2.07	0.54
2:CC:137:VAL:HA	2:CC:148:ILE:HD13	1.90	0.54
2:CC:42:LEU:O	2:CC:46:LEU:HB2	2.07	0.54
3:CD:192:ALA:HB3	3:CD:194:ILE:HG22	1.90	0.54
3:CD:96:ARG:HB3	3:CD:98:ASP:OD2	2.07	0.54
9:CJ:49:PHE:HE1	9:CJ:67:ILE:HD12	1.72	0.54
1:CA:1228:C:P	12:CM:113:LYS:HB2	2.47	0.54
21:CN:50:LEU:N	21:CN:51:PRO:HD2	2.16	0.54
17:CT:27:MET:O	17:CT:31:ILE:HG13	2.08	0.54
36:D2:30:VAL:HA	36:D2:33:ARG:HD3	1.90	0.54
23:DB:1355:G:O2'	23:DB:1356:G:H5'	2.08	0.54
23:DB:1742:U:H2'	23:DB:1743:G:C8	2.43	0.54
23:DB:1917:U:C2'	23:DB:1918:A:H5'	2.37	0.54
23:DB:1946:U:H2'	23:DB:1947:C:H6	1.72	0.54
23:DB:2327:A:H2'	23:DB:2328:A:C8	2.43	0.54
23:DB:2630:G:H2'	23:DB:2631:G:H8	1.71	0.54
23:DB:85:G:N2	23:DB:98:G:H1'	2.23	0.54
25:DC:255:LYS:C	25:DC:257:ARG:H	2.10	0.54
26:DD:107:VAL:H	26:DD:205:PRO:HA	1.71	0.54
47:DF:102:LEU:CD1	47:DF:103:ILE:HG13	2.37	0.54
48:DG:93:TYR:C	48:DG:94:ARG:HG3	2.27	0.54
40:DH:135:HIS:CG	40:DH:136:SER:N	2.76	0.54
27:DK:37:ASP:O	27:DK:61:VAL:HA	2.06	0.54
38:DM:19:GLY:H	38:DM:38:ARG:HH12	1.54	0.54
38:DM:10:ARG:HH11	38:DM:89:VAL:HG22	1.72	0.54
43:DO:72:ALA:HB2	43:DO:105:ALA:HB1	1.90	0.54
49:DR:62:GLU:O	49:DR:96:VAL:HA	2.06	0.54
46:DU:84:PHE:O	46:DU:85:ARG:HB2	2.07	0.54
46:DU:86:PHE:CD1	46:DU:90:LYS:HB2	2.42	0.54
52:DW:24:ARG:HD3	52:DW:65:LYS:HG2	1.89	0.54
1:AA:1032:G:N3	1:AA:1032:G:H3'	2.23	0.54
1:AA:1186:G:H4'	8:AI:111:GLU:OE1	2.08	0.54
1:AA:266:G:O2'	1:AA:267:C:H3'	2.07	0.54
1:AA:284:C:O2'	1:AA:285:C:H5'	2.07	0.54
1:AA:384:G:H2'	1:AA:385:C:H6	1.73	0.54
1:AA:601:G:H2'	1:AA:602:A:C8	2.43	0.54
1:AA:72:A:H2'	1:AA:73:C:C6	2.43	0.54
1:AA:878:A:H5''	7:AH:80:PRO:HG2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:129:VAL:HG12	3:AD:131:ILE:H	1.72	0.54
4:AE:87:VAL:HG23	4:AE:91:SER:O	2.07	0.54
6:AG:74:VAL:HG12	6:AG:87:PRO:HD3	1.90	0.54
7:AH:90:GLU:O	7:AH:91:LEU:C	2.46	0.54
8:AI:94:ARG:HA	8:AI:97:LEU:HD12	1.90	0.54
6:AG:149:ALA:HB2	10:AK:60:PHE:CB	2.37	0.54
22:BA:113:C:H2'	22:BA:114:C:C6	2.43	0.54
23:BB:1013:C:H2'	23:BB:1014:A:C8	2.42	0.54
23:BB:1060:U:C5	24:BI:131:THR:HG22	2.43	0.54
23:BB:143:C:H2'	23:BB:144:A:C8	2.43	0.54
23:BB:2087:G:H2'	23:BB:2088:A:C8	2.43	0.54
23:BB:2452:C:H2'	23:BB:2453:A:C8	2.43	0.54
23:BB:2457:U:O2'	23:BB:2458:G:H5'	2.07	0.54
23:BB:2462:C:H2'	23:BB:2463:C:H6	1.73	0.54
23:BB:2813:A:H2'	23:BB:2814:A:C8	2.43	0.54
23:BB:2828:G:H2'	23:BB:2829:A:H8	1.73	0.54
23:BB:2893:A:H4'	23:BB:2894:G:O5'	2.08	0.54
47:BF:35:LEU:HA	47:BF:153:ILE:HG12	1.89	0.54
23:BB:2746:U:H4'	48:BG:137:LYS:HB2	1.90	0.54
40:BH:131:SER:HB3	40:BH:140:ALA:C	2.28	0.54
40:BH:3:VAL:HA	40:BH:37:VAL:O	2.08	0.54
40:BH:63:ALA:HA	40:BH:66:ASN:CG	2.27	0.54
27:BK:2:ILE:HG13	27:BK:33:ALA:O	2.08	0.54
28:BP:8:GLU:HB3	28:BP:54:LEU:HB2	1.88	0.54
44:BQ:104:ALA:HA	49:BR:46:GLU:CD	2.28	0.54
49:BR:19:THR:HB	49:BR:97:LYS:HG3	1.90	0.54
1:CA:167:A:O2'	1:CA:168:G:H5'	2.08	0.54
1:CA:476:U:H2'	1:CA:477:C:H6	1.72	0.54
1:CA:513:C:H2'	1:CA:514:C:C6	2.43	0.54
3:CD:3:TYR:C	3:CD:4:LEU:HD12	2.28	0.54
4:CE:96:GLN:HB3	4:CE:123:LEU:HD12	1.89	0.54
6:CG:30:MET:CE	6:CG:33:GLY:HA2	2.38	0.54
10:CK:111:ASP:HB3	19:CU:19:LYS:HE2	1.89	0.54
15:CR:33:THR:HG23	15:CR:37:LYS:O	2.08	0.54
16:CS:40:PHE:HD2	16:CS:42:ASN:HD21	1.55	0.54
23:DB:117:G:H5'	23:DB:126:A:C8	2.40	0.54
23:DB:2542:A:H4'	23:DB:2543:G:H5'	1.89	0.54
23:DB:277:G:N2	23:DB:360:U:H3	2.06	0.54
23:DB:3:U:H2'	23:DB:4:U:H6	1.69	0.54
23:DB:643:A:C5	23:DB:644:A:N7	2.76	0.54
23:DB:672:C:O2'	23:DB:673:C:H5'	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:862:G:H2'	23:DB:863:A:C8	2.43	0.54
23:DB:934:U:H2'	23:DB:935:C:C6	2.43	0.54
27:DK:109:SER:C	27:DK:111:LYS:H	2.11	0.54
27:DK:2:ILE:HG13	27:DK:33:ALA:O	2.08	0.54
27:DK:19:VAL:HB	27:DK:41:ILE:HD11	1.90	0.54
52:DW:39:GLN:CG	52:DW:40:ARG:N	2.69	0.54
30:DY:6:ILE:O	30:DY:34:THR:HA	2.08	0.54
1:AA:1081:A:OP1	4:AE:22:LYS:HB2	2.08	0.54
5:AF:91:ARG:H	5:AF:93:LYS:HZ1	1.54	0.54
9:AJ:26:VAL:O	9:AJ:30:LYS:HG3	2.07	0.54
10:AK:28:ASN:HD21	10:AK:46:ALA:HB3	1.71	0.54
16:AS:10:ILE:HG22	16:AS:37:SER:CB	2.38	0.54
36:B2:18:PHE:O	36:B2:22:MET:HB2	2.08	0.54
23:BB:1163:G:O2'	23:BB:1164:C:H5'	2.07	0.54
23:BB:1374:G:O2'	23:BB:1375:U:H5'	2.07	0.54
23:BB:1758:U:O4'	23:BB:1758:U:O2	2.24	0.54
23:BB:2231:U:H2'	23:BB:2232:C:C6	2.42	0.54
23:BB:917:A:H5''	23:BB:2268:A:H61	1.73	0.54
23:BB:2848:G:N7	28:BP:94:ALA:HB2	2.22	0.54
23:BB:759:G:H2'	23:BB:760:G:H8	1.72	0.54
29:BE:3:LEU:HB2	29:BE:12:LEU:HB3	1.89	0.54
29:BE:48:THR:O	29:BE:52:VAL:HG23	2.08	0.54
41:BJ:93:ILE:O	41:BJ:97:PRO:HG3	2.07	0.54
28:BP:24:THR:OG1	28:BP:86:LYS:HB3	2.07	0.54
23:BB:141:G:N1	50:BT:1:MET:HA	2.23	0.54
52:BW:49:ASN:C	52:BW:50:VAL:HG22	2.28	0.54
39:BX:41:HIS:O	39:BX:44:LYS:HB3	2.08	0.54
1:CA:1148:U:H2'	1:CA:1149:C:O4'	2.08	0.54
1:CA:1496:C:H4'	23:DB:1920:C:O2'	2.07	0.54
1:CA:268:U:H2'	1:CA:269:C:C6	2.41	0.54
1:CA:89:U:H2'	1:CA:90:C:H6	1.73	0.54
18:CB:163:ILE:HG23	18:CB:164:ASP:N	2.17	0.54
5:CF:18:VAL:O	5:CF:22:ILE:HG13	2.08	0.54
6:CG:98:LEU:HD22	6:CG:102:TRP:NE1	2.23	0.54
1:CA:1124:G:H5'	9:CJ:37:ARG:HE	1.72	0.54
11:CL:98:ARG:HH21	11:CL:104:SER:C	2.10	0.54
16:CS:39:ILE:HD11	16:CS:68:HIS:CB	2.32	0.54
23:DB:1300:G:H4'	23:DB:1301:A:H5'	1.90	0.54
23:DB:1387:A:H2'	23:DB:1388:G:H8	1.73	0.54
23:DB:1692:U:H2'	23:DB:1694:C:C5	2.43	0.54
23:DB:1870:C:H2'	23:DB:1871:A:C8	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1130:U:C2	23:DB:2025:C:H5''	2.43	0.54
23:DB:2269:G:H4'	52:DW:19:ARG:HH11	1.73	0.54
23:DB:2580:U:C5'	26:DD:136:ASN:H	2.21	0.54
23:DB:2646:C:H2'	23:DB:2647:U:O4'	2.07	0.54
23:DB:423:A:H5''	23:DB:424:G:H5'	1.88	0.54
23:DB:441:U:H2'	23:DB:442:G:C8	2.43	0.54
23:DB:743:A:O2'	23:DB:744:U:H5'	2.08	0.54
23:DB:920:A:H2'	23:DB:921:C:H6	1.72	0.54
23:DB:967:U:H2'	23:DB:968:C:C6	2.43	0.54
25:DC:209:ALA:HA	25:DC:212:TRP:CE2	2.43	0.54
26:DD:125:TRP:CD2	26:DD:160:LYS:HB3	2.42	0.54
40:DH:94:ILE:HG23	40:DH:98:ASP:HB3	1.89	0.54
45:DS:42:LYS:O	45:DS:45:VAL:HG22	2.06	0.54
1:AA:1316:G:N1	1:AA:1318:A:H5''	2.23	0.54
1:AA:1332:A:H2'	1:AA:1333:A:C8	2.42	0.54
1:AA:255:G:H2'	1:AA:256:U:C6	2.42	0.54
3:AD:96:ARG:HB3	3:AD:98:ASP:OD2	2.08	0.54
4:AE:83:PRO:HB3	4:AE:96:GLN:HG3	1.90	0.54
6:AG:94:ARG:NH1	6:AG:98:LEU:HD21	2.23	0.54
8:AI:40:ARG:HA	8:AI:44:ARG:HD3	1.89	0.54
8:AI:98:ARG:C	8:AI:100:ALA:H	2.11	0.54
1:AA:1226:C:H5''	12:AM:94:LEU:HD11	1.90	0.54
22:BA:76:G:H2'	22:BA:77:U:H6	1.71	0.54
23:BB:1038:G:H2'	23:BB:1039:A:H8	1.73	0.54
23:BB:1300:G:H4'	23:BB:1301:A:H5'	1.89	0.54
23:BB:1459:G:H3'	23:BB:1460:U:C5'	2.38	0.54
23:BB:1571:A:H2'	23:BB:1572:A:C8	2.43	0.54
23:BB:2649:C:H2'	23:BB:2650:U:H6	1.71	0.54
23:BB:598:U:H2'	23:BB:599:A:C8	2.43	0.54
25:BC:138:SER:O	25:BC:162:GLN:HA	2.08	0.54
23:BB:2821:A:OP1	26:BD:115:GLY:HA3	2.07	0.54
47:BF:100:GLU:HA	47:BF:100:GLU:OE1	2.07	0.54
47:BF:102:LEU:HA	47:BF:106:ALA:CB	2.36	0.54
47:BF:110:ILE:CG2	47:BF:113:PHE:HB3	2.38	0.54
47:BF:57:ALA:HB2	47:BF:64:PRO:HG2	1.89	0.54
47:BF:95:MET:O	47:BF:99:PHE:HB2	2.08	0.54
24:BI:112:LYS:O	24:BI:116:MET:HG3	2.08	0.54
28:BP:50:ARG:CB	28:BP:56:SER:HB3	2.38	0.54
44:BQ:101:ASP:OD1	44:BQ:104:ALA:HB2	2.07	0.54
50:BT:35:ALA:O	50:BT:81:LYS:HB3	2.08	0.54
1:CA:1074:G:H2'	1:CA:1075:U:C6	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1250:A:H2'	1:CA:1251:A:H8	1.73	0.54
1:CA:50:A:H1'	1:CA:52:C:C6	2.43	0.54
18:CB:88:GLN:HE22	18:CB:217:ALA:HB1	1.72	0.54
18:CB:45:THR:HA	18:CB:48:MET:HG3	1.89	0.54
3:CD:13:ARG:HG3	3:CD:55:ARG:NH1	2.22	0.54
5:CF:69:GLU:O	5:CF:72:ASP:HB3	2.07	0.54
7:CH:94:VAL:HG21	7:CH:100:ILE:O	2.08	0.54
36:D2:16:HIS:HB3	36:D2:21:ARG:NH1	2.23	0.54
23:DB:1789:A:H2'	23:DB:1790:C:O4'	2.08	0.54
23:DB:1830:C:H2'	23:DB:1831:G:H8	1.73	0.54
23:DB:2104:C:H3'	23:DB:2104:C:C6	2.43	0.54
23:DB:2617:U:O2'	23:DB:2618:G:H5'	2.08	0.54
23:DB:2649:C:H2'	23:DB:2650:U:H6	1.72	0.54
23:DB:2714:G:O2'	23:DB:2715:C:H5'	2.08	0.54
23:DB:30:G:H2'	23:DB:31:C:H6	1.69	0.54
26:DD:5:VAL:HG23	26:DD:32:ASN:ND2	2.23	0.54
29:DE:34:ALA:CB	29:DE:96:VAL:HG21	2.38	0.54
40:DH:110:VAL:HG23	40:DH:132:PHE:CD2	2.43	0.54
27:DK:11:ALA:HB3	27:DK:85:VAL:HG22	1.88	0.54
37:DL:110:VAL:HB	37:DL:127:VAL:HG23	1.90	0.54
43:DO:35:ILE:HG13	43:DO:71:ALA:HB1	1.90	0.54
49:DR:49:ILE:HG12	49:DR:53:PHE:CA	2.38	0.54
46:DU:26:ASN:ND2	46:DU:26:ASN:N	2.55	0.54
39:DX:23:ARG:HD2	39:DX:27:ASN:ND2	2.22	0.54
1:AA:1418:A:N6	1:AA:1482:G:H1'	2.23	0.54
1:AA:1511:G:O2'	1:AA:1512:U:H5'	2.07	0.54
1:AA:235:C:H2'	1:AA:236:A:H8	1.71	0.54
1:AA:468:A:H3'	1:AA:469:C:H6	1.72	0.54
1:AA:476:U:O2'	1:AA:477:C:H5'	2.07	0.54
1:AA:489:C:H2'	1:AA:490:C:H6	1.73	0.54
2:AC:131:ARG:HH11	2:AC:135:ARG:HH12	1.55	0.54
6:AG:58:LEU:H	6:AG:58:LEU:CD2	2.18	0.54
7:AH:50:VAL:O	7:AH:50:VAL:HG22	2.08	0.54
7:AH:55:LYS:HE3	7:AH:56:PRO:HD2	1.90	0.54
16:AS:44:ILE:HG13	16:AS:62:THR:CA	2.36	0.54
23:BB:1340:U:O4	50:BT:64:LYS:HD2	2.08	0.54
23:BB:1447:C:H2'	23:BB:1448:G:C8	2.43	0.54
23:BB:1488:C:O2'	23:BB:1489:C:H5'	2.08	0.54
23:BB:1851:U:H2'	23:BB:1852:U:H6	1.72	0.54
23:BB:1854:A:H2	23:BB:2087:G:N3	2.05	0.54
23:BB:2298:A:H2'	23:BB:2299:U:O4'	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2438:U:H2'	23:BB:2441:U:H5	1.73	0.54
23:BB:2676:C:H2'	23:BB:2677:G:H8	1.73	0.54
23:BB:2722:G:H4'	42:BN:4:ARG:HB2	1.89	0.54
23:BB:2880:C:O2'	23:BB:2881:U:H5'	2.08	0.54
23:BB:743:A:O2'	23:BB:744:U:H5'	2.08	0.54
23:BB:925:A:O2'	23:BB:926:G:H5'	2.08	0.54
23:BB:93:G:H2'	23:BB:94:A:O4'	2.07	0.54
23:BB:2680:U:OP2	26:BD:114:LYS:HB3	2.08	0.54
23:BB:1007:C:O3'	41:BJ:110:PRO:HB3	2.07	0.54
37:BL:57:LEU:O	37:BL:61:LEU:HD13	2.08	0.54
43:BO:35:ILE:HG13	43:BO:71:ALA:HB1	1.90	0.54
46:BU:64:ILE:HD11	46:BU:68:ASN:ND2	2.23	0.54
46:BU:84:PHE:O	46:BU:85:ARG:HB2	2.08	0.54
35:BV:63:ILE:HD11	35:BV:72:VAL:HG22	1.88	0.54
1:CA:1070:U:H2'	1:CA:1071:C:C6	2.43	0.54
1:CA:1493:A:O5'	1:CA:1493:A:N3	2.41	0.54
1:CA:618:C:N3	1:CA:622:A:N6	2.53	0.54
1:CA:772:U:H2'	1:CA:773:G:O4'	2.07	0.54
18:CB:130:LYS:O	18:CB:134:LEU:HD23	2.07	0.54
18:CB:114:LYS:HE2	18:CB:151:LYS:HB2	1.90	0.54
18:CB:218:ALA:HA	18:CB:221:ARG:CZ	2.38	0.54
4:CE:155:LYS:HD2	7:CH:70:VAL:O	2.07	0.54
20:CO:8:THR:O	20:CO:12:VAL:HG23	2.07	0.54
16:CS:43:MET:HB2	16:CS:61:VAL:HG21	1.90	0.54
23:DB:2199:A:H3'	23:DB:2200:C:H6	1.72	0.54
23:DB:2298:A:H2'	23:DB:2299:U:O4'	2.08	0.54
23:DB:2869:G:H2'	23:DB:2870:C:H6	1.72	0.54
23:DB:692:C:H2'	23:DB:693:A:H8	1.73	0.54
23:DB:951:C:O2'	23:DB:952:G:H5'	2.08	0.54
23:DB:2772:C:H4'	26:DD:171:THR:HG21	1.89	0.54
48:DG:67:ALA:O	48:DG:71:LEU:HD23	2.08	0.54
40:DH:39:ALA:C	40:DH:41:LYS:H	2.12	0.54
27:DK:25:LEU:HD21	27:DK:40:LYS:H	1.71	0.54
37:DL:142:ILE:N	37:DL:142:ILE:HD12	2.23	0.54
42:DN:24:MET:SD	42:DN:44:LEU:HD22	2.48	0.54
1:AA:1225:A:H5''	1:AA:1226:C:C5	2.38	0.54
1:AA:1290:G:H2'	1:AA:1291:U:H6	1.72	0.54
1:AA:1396:A:C4'	1:AA:1397:C:H5'	2.38	0.54
1:AA:1484:C:H2'	1:AA:1485:U:H6	1.72	0.54
1:AA:413:G:O6	3:AD:32:LYS:HE2	2.07	0.54
18:AB:98:GLY:C	18:AB:100:LEU:H	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:213:LEU:O	18:AB:216:VAL:HG22	2.07	0.54
5:AF:84:VAL:HG22	5:AF:85:ILE:N	2.22	0.54
9:AJ:7:ARG:HG3	9:AJ:101:SER:HB2	1.90	0.54
12:AM:1:ALA:O	12:AM:3:ILE:HG13	2.07	0.54
12:AM:44:ILE:CA	12:AM:47:LEU:HD23	2.36	0.54
23:BB:1584:U:H5'	23:BB:1585:C:C5	2.44	0.54
23:BB:1593:A:H2'	23:BB:1594:U:O4'	2.08	0.54
23:BB:2322:A:H3'	23:BB:2323:G:H8	1.73	0.54
23:BB:2547:A:H5'	23:BB:2566:A:C2	2.43	0.54
23:BB:2605:U:H2'	23:BB:2606:C:C6	2.42	0.54
23:BB:2861:U:H2'	23:BB:2862:G:C8	2.43	0.54
23:BB:581:C:H2'	23:BB:582:A:C8	2.43	0.54
25:BC:66:PHE:HB2	25:BC:150:GLY:O	2.08	0.54
47:BF:102:LEU:CD1	47:BF:103:ILE:HG13	2.37	0.54
40:BH:116:ARG:NE	40:BH:131:SER:HB2	2.22	0.54
27:BK:115:ILE:HG23	27:BK:116:ILE:N	2.22	0.54
38:BM:34:LYS:HA	38:BM:100:LYS:O	2.08	0.54
43:BO:52:SER:O	43:BO:58:ILE:HD12	2.08	0.54
49:BR:60:LYS:H	49:BR:100:GLY:HA3	1.72	0.54
50:BT:36:LYS:O	50:BT:36:LYS:HD3	2.08	0.54
52:BW:38:ARG:N	52:BW:56:HIS:HD2	2.06	0.54
1:CA:1343:G:H2'	1:CA:1344:C:H6	1.72	0.54
2:CC:146:LYS:HB2	2:CC:202:PHE:HE2	1.71	0.54
8:CI:11:ARG:HA	8:CI:105:ARG:NH1	2.23	0.54
8:CI:61:ASP:O	8:CI:62:LEU:HD13	2.07	0.54
10:CK:70:ALA:HA	10:CK:73:VAL:CG2	2.38	0.54
11:CL:107:LYS:C	11:CL:109:ARG:H	2.11	0.54
11:CL:14:LYS:HG2	11:CL:15:VAL:N	2.22	0.54
22:DA:78:A:H2'	22:DA:79:G:O4'	2.07	0.54
23:DB:1470:A:H2'	23:DB:1471:G:O4'	2.08	0.54
23:DB:1704:C:H2'	23:DB:1705:A:H8	1.72	0.54
23:DB:979:A:H3'	23:DB:980:A:C5'	2.38	0.54
25:DC:245:THR:O	25:DC:247:TRP:N	2.41	0.54
26:DD:46:ARG:NH1	26:DD:86:GLU:H	2.06	0.54
47:DF:2:LYS:HB2	47:DF:100:GLU:OE2	2.07	0.54
40:DH:131:SER:HB2	40:DH:140:ALA:O	2.07	0.54
40:DH:4:ILE:HG23	40:DH:16:GLY:HA2	1.90	0.54
23:DB:1099:G:H4'	24:DI:4:VAL:CG1	2.38	0.54
37:DL:123:ARG:HA	37:DL:143:GLU:CB	2.31	0.54
37:DL:110:VAL:HG23	37:DL:126:ARG:O	2.08	0.54
37:DL:19:LEU:HD23	37:DL:27:LEU:HD23	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1190:G:H5''	37:DL:32:GLY:O	2.07	0.54
37:DL:4:ASN:HD22	37:DL:4:ASN:N	2.06	0.54
44:DQ:90:ASP:H	49:DR:39:LEU:CD1	2.21	0.54
51:DZ:27:ARG:HG3	51:DZ:28:ARG:H	1.72	0.54
1:AA:1238:A:H2'	1:AA:1239:A:H8	1.72	0.53
1:AA:205:A:H2'	1:AA:206:C:C6	2.42	0.53
1:AA:253:A:H2'	1:AA:254:G:C8	2.44	0.53
1:AA:87:C:H2'	1:AA:88:U:C4'	2.37	0.53
18:AB:68:PHE:HE1	18:AB:88:GLN:HB3	1.73	0.53
4:AE:95:MET:HA	4:AE:124:ALA:CB	2.38	0.53
5:AF:69:GLU:O	5:AF:73:GLU:HG2	2.09	0.53
6:AG:129:ASN:HA	6:AG:134:VAL:HG21	1.90	0.53
6:AG:16:LYS:HB3	6:AG:43:TYR:CE1	2.42	0.53
10:AK:80:ASN:CA	10:AK:105:ARG:HB3	2.38	0.53
11:AL:107:LYS:C	11:AL:109:ARG:H	2.11	0.53
20:AO:25:THR:HG21	20:AO:70:LEU:HD23	1.90	0.53
19:AU:39:LYS:N	19:AU:40:PRO:CD	2.71	0.53
36:B2:3:ARG:HA	36:B2:3:ARG:CZ	2.38	0.53
36:B2:3:ARG:HA	36:B2:3:ARG:NE	2.22	0.53
23:BB:1015:U:H2'	23:BB:1016:G:C8	2.42	0.53
23:BB:1742:U:H2'	23:BB:1743:G:C8	2.43	0.53
23:BB:1771:C:H2'	23:BB:1772:A:H8	1.73	0.53
23:BB:397:U:OP1	51:BZ:31:PRO:HA	2.07	0.53
23:BB:423:A:H5''	23:BB:424:G:H5'	1.90	0.53
23:BB:526:A:N6	23:BB:2626:C:H4'	2.23	0.53
23:BB:736:C:H2'	23:BB:737:C:C6	2.43	0.53
29:BE:146:VAL:HG11	29:BE:187:VAL:HG23	1.90	0.53
29:BE:47:LYS:HB3	29:BE:51:GLU:HB2	1.90	0.53
1:CA:1164:G:H2'	1:CA:1165:U:C6	2.43	0.53
1:CA:1280:A:O4'	9:CJ:43:PRO:HG3	2.07	0.53
1:CA:1469:C:H2'	1:CA:1470:U:O4'	2.07	0.53
1:CA:621:A:H2'	1:CA:622:A:H8	1.72	0.53
1:CA:882:C:O2'	1:CA:883:C:H5'	2.08	0.53
1:CA:923:A:H2'	1:CA:924:C:H6	1.73	0.53
4:CE:92:ARG:HB3	4:CE:92:ARG:NH1	2.23	0.53
4:CE:98:ALA:HB2	4:CE:123:LEU:HG	1.91	0.53
8:CI:113:LYS:HA	8:CI:120:ALA:HB2	1.90	0.53
12:CM:33:LEU:CD1	12:CM:40:GLU:HA	2.38	0.53
21:CN:3:GLN:HA	21:CN:6:LYS:CG	2.38	0.53
23:DB:2073:C:O2'	23:DB:2074:U:H5'	2.07	0.53
23:DB:208:C:H2'	23:DB:209:C:H6	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:67:U:H2'	23:DB:68:G:C8	2.41	0.53
23:DB:93:G:H2'	23:DB:94:A:O4'	2.08	0.53
25:DC:138:SER:O	25:DC:162:GLN:HA	2.08	0.53
29:DE:3:LEU:HB2	29:DE:12:LEU:HB3	1.90	0.53
29:DE:2:GLU:HA	29:DE:13:THR:CA	2.35	0.53
29:DE:29:HIS:O	29:DE:33:VAL:HG23	2.09	0.53
12:CM:66:GLY:HA2	47:DF:143:ASP:HB3	1.88	0.53
38:DM:21:ALA:HB3	38:DM:99:GLY:O	2.09	0.53
28:DP:20:ARG:NH2	28:DP:91:VAL:HG22	2.23	0.53
44:DQ:43:GLN:NE2	49:DR:77:PHE:HB3	2.23	0.53
44:DQ:63:ARG:HH12	44:DQ:96:ASP:CB	2.13	0.53
39:DX:39:GLN:O	39:DX:42:LEU:HB2	2.08	0.53
23:DB:96:C:H4'	39:DX:41:HIS:CE1	2.43	0.53
1:AA:1118:U:H1'	1:AA:1179:A:C4	2.43	0.53
1:AA:1336:C:H4'	1:AA:1337:G:C4	2.43	0.53
1:AA:1439:G:H2'	1:AA:1440:U:O4'	2.07	0.53
1:AA:493:A:H3'	1:AA:494:G:C8	2.44	0.53
1:AA:570:G:H2'	1:AA:571:U:C6	2.43	0.53
18:AB:10:LYS:O	18:AB:13:VAL:HG22	2.07	0.53
4:AE:101:GLY:H	4:AE:121:ASN:ND2	2.05	0.53
6:AG:125:ASP:HA	6:AG:128:GLU:OE1	2.08	0.53
6:AG:94:ARG:HD3	6:AG:98:LEU:HD11	1.89	0.53
8:AI:40:ARG:N	8:AI:44:ARG:HD3	2.23	0.53
34:B3:18:LYS:HD2	34:B3:19:GLY:N	2.23	0.53
23:BB:1268:A:H2'	23:BB:1269:A:O4'	2.08	0.53
23:BB:1387:A:H2'	23:BB:1388:G:H8	1.73	0.53
23:BB:1923:U:H2'	23:BB:1924:C:C6	2.43	0.53
23:BB:2256:G:H2'	23:BB:2257:U:C6	2.43	0.53
23:BB:2857:G:N2	23:BB:2859:G:H3'	2.23	0.53
23:BB:594:U:H2'	23:BB:595:C:C6	2.43	0.53
23:BB:814:C:O2'	23:BB:815:C:H5'	2.07	0.53
25:BC:75:ALA:CB	25:BC:95:TYR:HA	2.39	0.53
48:BG:76:ILE:HB	48:BG:82:PHE:CZ	2.44	0.53
40:BH:99:ILE:CD1	40:BH:144:VAL:HB	2.37	0.53
41:BJ:44:TYR:O	41:BJ:45:THR:HB	2.09	0.53
38:BM:64:TRP:HB2	38:BM:104:GLU:HB2	1.90	0.53
49:BR:14:VAL:HG22	49:BR:15:SER:N	2.23	0.53
44:BQ:111:LYS:HE2	49:BR:50:GLY:HA3	1.89	0.53
35:BV:6:ALA:HB3	35:BV:65:VAL:HG12	1.90	0.53
30:BY:2:LYS:HD3	30:BY:2:LYS:H	1.74	0.53
1:CA:126:G:H4'	1:CA:634:C:H1'	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1365:G:H2'	1:CA:1366:C:C6	2.42	0.53
1:CA:648:A:H2'	1:CA:649:A:C8	2.43	0.53
1:CA:861:G:H2'	1:CA:862:C:H6	1.73	0.53
18:CB:53:LEU:HD11	18:CB:216:VAL:HA	1.90	0.53
18:CB:66:ILE:HD12	18:CB:159:ALA:HB3	1.88	0.53
2:CC:71:ARG:HB2	2:CC:74:ILE:HG21	1.89	0.53
3:CD:129:VAL:HG12	3:CD:131:ILE:H	1.73	0.53
5:CF:3:HIS:ND1	5:CF:95:ALA:N	2.56	0.53
6:CG:134:VAL:HG12	6:CG:137:ARG:HH21	1.73	0.53
11:CL:17:LYS:HD2	11:CL:18:SER:O	2.09	0.53
12:CM:65:GLU:N	12:CM:68:LEU:HD12	2.23	0.53
20:CO:64:ARG:CA	20:CO:64:ARG:HE	2.19	0.53
36:D2:18:PHE:O	36:D2:22:MET:HB2	2.08	0.53
34:D3:40:LYS:O	34:D3:43:LEU:HB2	2.09	0.53
22:DA:54:G:O2'	22:DA:55:U:H5'	2.08	0.53
23:DB:2028:U:H2'	23:DB:2029:G:C8	2.43	0.53
23:DB:2313:C:H4'	47:DF:87:LYS:HB3	1.90	0.53
23:DB:480:A:H2'	23:DB:480:A:N3	2.24	0.53
26:DD:3:GLY:C	26:DD:4:LEU:HD22	2.29	0.53
47:DF:41:GLU:HG3	47:DF:48:LEU:HD11	1.89	0.53
48:DG:83:THR:C	48:DG:84:LYS:HD3	2.29	0.53
48:DG:84:LYS:CG	48:DG:85:LYS:H	2.20	0.53
27:DK:34:GLY:O	27:DK:36:GLY:N	2.42	0.53
42:DN:62:ASN:HD22	42:DN:62:ASN:N	2.07	0.53
43:DO:111:ARG:HA	43:DO:115:LEU:O	2.08	0.53
43:DO:76:LYS:O	43:DO:80:GLU:HG2	2.08	0.53
28:DP:50:ARG:CB	28:DP:56:SER:HB3	2.38	0.53
50:DT:36:LYS:O	50:DT:36:LYS:HD3	2.08	0.53
30:DY:7:THR:HG22	30:DY:8:GLN:N	2.22	0.53
1:AA:1237:C:C3'	1:AA:1238:A:H5'	2.37	0.53
18:AB:177:ASN:O	18:AB:178:LEU:HD23	2.08	0.53
2:AC:31:ASN:ND2	2:AC:58:ARG:HH22	2.06	0.53
7:AH:28:SER:HB2	7:AH:56:PRO:HB2	1.89	0.53
4:AE:156:ARG:HH12	7:AH:42:GLU:HB3	1.70	0.53
20:AO:84:ARG:C	20:AO:85:LEU:HD12	2.28	0.53
36:B2:3:ARG:HE	36:B2:4:THR:HG22	1.73	0.53
22:BA:78:A:H2'	22:BA:79:G:O4'	2.08	0.53
23:BB:1230:A:H2'	23:BB:1231:U:H6	1.74	0.53
23:BB:1461:C:H2'	23:BB:1462:C:C6	2.41	0.53
23:BB:1729:U:H5	23:BB:1731:G:N2	2.06	0.53
23:BB:528:A:N1	23:BB:2042:A:H2'	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2080:A:OP1	51:BZ:20:HIS:HB3	2.09	0.53
23:BB:2645:G:H4'	23:BB:2732:G:H2'	1.91	0.53
37:BL:100:ILE:O	37:BL:100:ILE:HG12	2.08	0.53
42:BN:49:GLU:HB2	42:BN:50:PRO:HD3	1.90	0.53
42:BN:77:ALA:O	42:BN:81:ASN:HB2	2.08	0.53
50:BT:50:LEU:N	50:BT:50:LEU:HD22	2.22	0.53
46:BU:98:ASN:OD1	46:BU:100:GLU:HB2	2.08	0.53
51:BZ:31:PRO:HB2	51:BZ:33:LEU:HG	1.90	0.53
1:CA:195:A:H1'	1:CA:222:C:O2'	2.08	0.53
1:CA:336:A:O2'	1:CA:337:G:H5'	2.08	0.53
1:CA:468:A:H3'	1:CA:469:C:H6	1.72	0.53
1:CA:600:A:H2'	1:CA:601:G:H8	1.72	0.53
1:CA:678:U:H2'	1:CA:679:C:C6	2.43	0.53
18:CB:75:ALA:O	18:CB:79:VAL:N	2.39	0.53
6:CG:98:LEU:HD22	6:CG:102:TRP:CE2	2.43	0.53
8:CI:83:THR:HG21	8:CI:102:PHE:O	2.08	0.53
8:CI:8:THR:HG1	8:CI:9:GLY:H	1.54	0.53
11:CL:3:VAL:O	11:CL:7:VAL:HG23	2.09	0.53
12:CM:19:THR:C	12:CM:21:ILE:H	2.12	0.53
9:CJ:51:VAL:HG22	21:CN:80:ARG:HG3	1.91	0.53
20:CO:47:LYS:C	20:CO:49:ASP:H	2.11	0.53
23:DB:1315:C:H2'	23:DB:1316:U:H6	1.73	0.53
23:DB:1376:C:H3'	56:DB:3278:HOH:O	2.08	0.53
23:DB:1392:A:H2'	23:DB:1393:A:C8	2.44	0.53
23:DB:1414:C:H2'	23:DB:1415:U:H6	1.73	0.53
23:DB:1637:A:H2'	23:DB:1638:C:C6	2.43	0.53
23:DB:1758:U:O4'	23:DB:1758:U:O2	2.24	0.53
23:DB:2194:U:H2'	23:DB:2195:U:C6	2.43	0.53
23:DB:2728:U:H2'	23:DB:2729:G:C8	2.43	0.53
23:DB:2901:C:O2'	23:DB:2902:C:H5'	2.09	0.53
25:DC:202:ARG:HE	25:DC:213:ARG:NH2	2.06	0.53
25:DC:20:ASN:O	25:DC:23:LEU:HB2	2.08	0.53
26:DD:24:VAL:HG21	26:DD:188:LEU:HB3	1.90	0.53
47:DF:65:LEU:N	47:DF:88:VAL:HG22	2.22	0.53
48:DG:34:ARG:HD3	48:DG:34:ARG:N	2.22	0.53
24:DI:2:LYS:C	24:DI:3:LYS:HD2	2.28	0.53
42:DN:77:ALA:O	42:DN:81:ASN:HB2	2.09	0.53
27:DK:75:SER:HA	28:DP:72:VAL:O	2.08	0.53
45:DS:68:ASP:C	45:DS:69:LEU:HD12	2.29	0.53
1:AA:1453:G:H2'	1:AA:1454:G:O4'	2.09	0.53
1:AA:926:G:H3'	1:AA:1505:G:H21	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:8:A:C6	3:AD:205:LYS:HB2	2.43	0.53
5:AF:69:GLU:O	5:AF:72:ASP:HB3	2.09	0.53
17:AT:79:THR:O	17:AT:82:ILE:HG13	2.08	0.53
22:BA:37:C:H2'	22:BA:38:C:O4'	2.08	0.53
23:BB:1425:G:H2'	23:BB:1426:G:C8	2.43	0.53
23:BB:2053:G:O2'	23:BB:2054:A:H5'	2.08	0.53
23:BB:279:A:C2	23:BB:362:A:H4'	2.44	0.53
25:BC:42:ARG:NH2	25:BC:48:ILE:HD11	2.24	0.53
26:BD:106:LYS:H	26:BD:106:LYS:HD3	1.73	0.53
26:BD:149:ASN:C	26:BD:152:PRO:HD2	2.29	0.53
47:BF:115:GLY:HA3	47:BF:177:ARG:HD2	1.89	0.53
47:BF:35:LEU:HD22	47:BF:56:LEU:HD11	1.89	0.53
48:BG:24:THR:HG22	48:BG:34:ARG:HA	1.90	0.53
24:BI:18:ASN:N	24:BI:19:PRO:CD	2.71	0.53
27:BK:37:ASP:O	27:BK:61:VAL:HA	2.08	0.53
38:BM:35:ALA:H	38:BM:100:LYS:H	1.55	0.53
38:BM:19:GLY:H	38:BM:38:ARG:HH12	1.55	0.53
38:BM:59:ARG:HH11	38:BM:60:GLN:HB3	1.73	0.53
38:BM:17:ASN:HD21	38:BM:95:LEU:HD11	1.74	0.53
45:BS:31:GLN:O	45:BS:34:ASP:HB2	2.08	0.53
30:BY:37:ARG:HG3	30:BY:38:GLU:OE1	2.09	0.53
1:CA:1250:A:H2'	1:CA:1251:A:C8	2.43	0.53
1:CA:34:C:H2'	1:CA:35:G:C8	2.44	0.53
1:CA:613:C:H2'	1:CA:614:C:C6	2.44	0.53
1:CA:77:A:H2'	1:CA:78:A:C8	2.43	0.53
1:CA:842:U:OP1	1:CA:843:U:H4'	2.08	0.53
2:CC:6:PRO:HG2	2:CC:183:TYR:CG	2.43	0.53
14:CQ:24:ILE:HD13	14:CQ:43:LEU:HD13	1.89	0.53
31:D0:38:LEU:HB3	31:D0:41:HIS:NE2	2.24	0.53
23:DB:1904:G:O2'	23:DB:1905:C:H5'	2.08	0.53
23:DB:198:C:H6	23:DB:198:C:O5'	1.92	0.53
23:DB:2817:U:O2'	23:DB:2837:A:H1'	2.09	0.53
23:DB:594:U:H2'	23:DB:595:C:C6	2.42	0.53
23:DB:799:G:H3'	23:DB:800:A:H2'	1.90	0.53
23:DB:90:U:H3'	23:DB:91:A:C5'	2.38	0.53
25:DC:129:LEU:HB3	25:DC:134:ILE:HG22	1.91	0.53
25:DC:75:ALA:CB	25:DC:95:TYR:HA	2.39	0.53
25:DC:91:ALA:HB2	25:DC:105:ALA:HB2	1.91	0.53
26:DD:125:TRP:CE2	26:DD:160:LYS:HB3	2.44	0.53
47:DF:8:LYS:HA	47:DF:12:VAL:CG2	2.36	0.53
40:DH:116:ARG:O	40:DH:130:VAL:HG13	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DH:50:ARG:C	40:DH:52:ALA:H	2.11	0.53
40:DH:53:GLU:C	40:DH:55:GLU:H	2.12	0.53
38:DM:41:LEU:HB3	38:DM:46:ILE:HG22	1.91	0.53
43:DO:115:LEU:C	43:DO:116:GLN:HE21	2.11	0.53
28:DP:99:LEU:O	28:DP:99:LEU:HD13	2.08	0.53
49:DR:19:THR:HB	49:DR:97:LYS:HG3	1.90	0.53
49:DR:5:PHE:HD1	49:DR:5:PHE:H	1.55	0.53
50:DT:39:THR:HG22	50:DT:42:GLU:HG2	1.90	0.53
1:AA:121:U:H4'	1:AA:122:G:N7	2.23	0.53
1:AA:1292:G:H2'	1:AA:1293:C:C6	2.43	0.53
1:AA:1499:A:H5'	1:AA:1499:A:C8	2.38	0.53
2:AC:3:LYS:HB2	2:AC:3:LYS:NZ	2.24	0.53
4:AE:81:GLN:H	4:AE:146:MET:CE	2.18	0.53
5:AF:55:HIS:HB2	5:AF:56:LYS:HE2	1.91	0.53
5:AF:80:PHE:CZ	25:BC:123:ILE:HG12	2.43	0.53
11:AL:14:LYS:HG2	11:AL:15:VAL:N	2.23	0.53
22:BA:55:U:H2'	22:BA:56:G:H8	1.74	0.53
23:BB:1470:A:H2'	23:BB:1471:G:O4'	2.07	0.53
23:BB:1599:U:H2'	23:BB:1600:C:C6	2.42	0.53
23:BB:2199:A:H3'	23:BB:2200:C:H6	1.73	0.53
23:BB:21:A:H2'	23:BB:22:C:C6	2.44	0.53
23:BB:237:C:O2'	23:BB:238:C:H5'	2.07	0.53
23:BB:2408:U:O2'	23:BB:2409:G:H5'	2.09	0.53
23:BB:2757:A:H2'	23:BB:2758:A:H5'	1.91	0.53
23:BB:2772:C:H4'	26:BD:171:THR:HG21	1.89	0.53
23:BB:915:C:H3'	23:BB:916:G:H8	1.72	0.53
47:BF:40:GLY:O	47:BF:41:GLU:HB2	2.08	0.53
38:BM:108:VAL:HG21	38:BM:112:LEU:HD12	1.89	0.53
38:BM:114:ARG:HG3	38:BM:130:PHE:CG	2.44	0.53
28:BP:20:ARG:NH2	28:BP:91:VAL:HG22	2.23	0.53
1:CA:1070:U:H2'	1:CA:1071:C:H6	1.72	0.53
1:CA:1171:A:H2'	1:CA:1172:C:C5	2.44	0.53
1:CA:778:G:H2'	1:CA:779:C:C6	2.44	0.53
1:CA:784:A:H2'	1:CA:785:G:C8	2.44	0.53
4:CE:106:ALA:HB1	4:CE:110:MET:HB3	1.89	0.53
4:CE:56:PRO:HG2	4:CE:57:ALA:H	1.73	0.53
4:CE:87:VAL:HG23	4:CE:91:SER:O	2.08	0.53
7:CH:90:GLU:O	7:CH:91:LEU:C	2.46	0.53
19:CU:42:THR:HB	19:CU:46:ARG:HE	1.73	0.53
33:D1:3:GLY:O	33:D1:5:ARG:N	2.42	0.53
23:DB:1098:A:C2'	24:DI:4:VAL:N	2.71	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1301:A:O2'	23:DB:1302:A:H3'	2.07	0.53
23:DB:1439:A:H1'	23:DB:1553:A:N6	2.23	0.53
23:DB:1560:G:H2'	23:DB:1561:C:C6	2.42	0.53
23:DB:959:A:O2'	23:DB:960:A:H5'	2.08	0.53
25:DC:152:GLN:C	25:DC:153:LEU:HD23	2.28	0.53
25:DC:202:ARG:HE	25:DC:213:ARG:HH21	1.56	0.53
48:DG:94:ARG:HH21	48:DG:104:LEU:HA	1.73	0.53
48:DG:3:VAL:O	48:DG:68:ARG:HG3	2.08	0.53
24:DI:57:VAL:HG23	24:DI:71:LYS:NZ	2.24	0.53
41:DJ:28:LEU:HG	41:DJ:29:ALA:N	2.22	0.53
41:DJ:44:TYR:O	41:DJ:45:THR:HB	2.08	0.53
27:DK:71:ARG:HA	27:DK:71:ARG:HE	1.74	0.53
51:DZ:31:PRO:HB2	51:DZ:33:LEU:HG	1.90	0.53
1:AA:1000:A:H2'	1:AA:1001:C:C6	2.43	0.53
1:AA:1014:A:H5''	16:AS:13:HIS:HB3	1.90	0.53
1:AA:1524:C:H2'	1:AA:1525:G:C8	2.43	0.53
1:AA:377:G:H2'	1:AA:378:G:C8	2.44	0.53
2:AC:137:VAL:HA	2:AC:148:ILE:HD13	1.90	0.53
2:AC:52:SER:CA	2:AC:113:LYS:HG2	2.38	0.53
2:AC:87:ARG:HE	2:AC:87:ARG:C	2.11	0.53
3:AD:55:ARG:NE	3:AD:55:ARG:HA	2.24	0.53
6:AG:74:VAL:HA	6:AG:86:VAL:O	2.08	0.53
19:AU:24:LYS:HD2	19:AU:25:ALA:N	2.24	0.53
33:B1:26:LYS:HB2	33:B1:52:LYS:HZ2	1.74	0.53
23:BB:143:C:N4	23:BB:144:A:N6	2.56	0.53
23:BB:1441:G:H2'	23:BB:1442:U:H6	1.71	0.53
23:BB:1830:C:H2'	23:BB:1831:G:H8	1.73	0.53
23:BB:871:U:H2'	23:BB:872:U:C6	2.44	0.53
25:BC:43:ASN:ND2	25:BC:44:ASN:H	2.05	0.53
26:BD:24:VAL:HG21	26:BD:188:LEU:HB3	1.91	0.53
29:BE:29:HIS:O	29:BE:33:VAL:HG23	2.09	0.53
47:BF:120:SER:HB2	47:BF:129:MET:HG3	1.90	0.53
47:BF:128:SER:HB3	47:BF:154:THR:HA	1.90	0.53
27:BK:39:ILE:O	27:BK:59:LYS:HA	2.09	0.53
37:BL:80:SER:H	37:BL:113:ALA:CB	2.21	0.53
38:BM:41:LEU:HB3	38:BM:46:ILE:HG22	1.90	0.53
43:BO:116:GLN:HE21	43:BO:116:GLN:N	2.06	0.53
44:BQ:105:PHE:HA	44:BQ:108:LEU:CD1	2.35	0.53
1:CA:1074:G:H2'	1:CA:1075:U:H6	1.72	0.53
1:CA:961:U:O4'	1:CA:961:U:O2	2.24	0.53
6:CG:66:GLU:HA	6:CG:69:ARG:HD2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CH:124:ILE:HG22	7:CH:125:ILE:N	2.23	0.53
1:CA:1348:U:H4'	8:CI:121:ARG:HG3	1.90	0.53
1:CA:1129:C:H5''	8:CI:17:ARG:HH12	1.74	0.53
9:CJ:44:THR:HG21	9:CJ:70:HIS:HD2	1.73	0.53
20:CO:46:HIS:O	20:CO:48:LYS:N	2.41	0.53
1:CA:958:A:P	16:CS:54:ARG:HH22	2.32	0.53
19:CU:39:LYS:N	19:CU:40:PRO:CD	2.71	0.53
32:D4:24:ARG:HG2	32:D4:36:ARG:CG	2.39	0.53
23:DB:118:A:OP2	23:DB:119:A:H5''	2.08	0.53
23:DB:1967:C:H2'	23:DB:1968:G:O4'	2.09	0.53
23:DB:2087:G:H2'	23:DB:2088:A:C8	2.44	0.53
23:DB:216:A:H2'	23:DB:217:A:H8	1.74	0.53
23:DB:2230:G:H2'	23:DB:2231:U:H6	1.72	0.53
23:DB:2857:G:N2	23:DB:2859:G:H3'	2.23	0.53
23:DB:2875:C:H2'	23:DB:2876:G:H8	1.73	0.53
23:DB:483:A:H5''	46:DU:46:LYS:CG	2.39	0.53
23:DB:608:A:H2'	23:DB:609:A:H8	1.74	0.53
23:DB:64:A:H2'	23:DB:65:U:C6	2.44	0.53
25:DC:247:TRP:O	25:DC:249:VAL:HG22	2.08	0.53
24:DI:52:LEU:HD22	24:DI:81:LYS:HD3	1.89	0.53
24:DI:99:LYS:HD3	24:DI:99:LYS:N	2.24	0.53
27:DK:116:ILE:HG13	27:DK:117:SER:N	2.23	0.53
23:DB:2700:A:H2	42:DN:71:ARG:HH12	1.56	0.53
49:DR:28:ALA:HB3	49:DR:31:GLU:HG3	1.90	0.53
50:DT:29:THR:HA	50:DT:86:THR:CA	2.34	0.53
1:AA:1044:A:C5	1:AA:1045:C:H1'	2.44	0.53
1:AA:268:U:H2'	1:AA:269:C:C6	2.44	0.53
1:AA:621:A:H2'	1:AA:622:A:H8	1.72	0.53
2:AC:131:ARG:HB3	2:AC:135:ARG:NH1	2.24	0.53
4:AE:158:LYS:NZ	7:AH:63:LYS:HD3	2.24	0.53
6:AG:84:TYR:O	6:AG:86:VAL:HG23	2.09	0.53
9:AJ:77:VAL:HB	9:AJ:78:GLU:OE1	2.09	0.53
11:AL:43:LYS:HE3	11:AL:44:PRO:HD3	1.90	0.53
11:AL:52:CYS:SG	11:AL:66:ILE:HD11	2.49	0.53
12:AM:37:GLY:O	12:AM:38:ILE:HD13	2.09	0.53
2:AC:25:THR:HG23	21:AN:75:LYS:HD2	1.90	0.53
32:B4:24:ARG:HG2	32:B4:36:ARG:CG	2.39	0.53
23:BB:1210:G:N3	23:BB:1212:G:N2	2.56	0.53
23:BB:1258:U:H2'	23:BB:1259:G:C8	2.43	0.53
23:BB:1789:A:H2'	23:BB:1790:C:O4'	2.09	0.53
23:BB:2651:C:O2'	23:BB:2652:C:H5'	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2674:G:H2'	23:BB:2675:A:C8	2.44	0.53
23:BB:408:G:O2'	23:BB:409:G:H5'	2.09	0.53
23:BB:644:A:H2'	23:BB:644:A:N3	2.24	0.53
23:BB:934:U:H2'	23:BB:935:C:H6	1.74	0.53
23:BB:951:C:O2'	23:BB:952:G:H5'	2.09	0.53
25:BC:4:LYS:CD	25:BC:5:CYS:H	2.22	0.53
26:BD:51:THR:HG22	26:BD:52:THR:H	1.73	0.53
29:BE:4:VAL:HG12	29:BE:5:LEU:N	2.21	0.53
29:BE:48:THR:C	29:BE:50:ALA:H	2.10	0.53
24:BI:81:LYS:HG3	24:BI:82:ALA:N	2.23	0.53
27:BK:99:ILE:HB	27:BK:118:LEU:HD22	1.90	0.53
43:BO:111:ARG:HA	43:BO:115:LEU:O	2.08	0.53
43:BO:28:VAL:O	43:BO:28:VAL:HG13	2.09	0.53
44:BQ:33:VAL:HG23	44:BQ:34:ALA:N	2.24	0.53
44:BQ:57:ARG:HA	44:BQ:60:TRP:CE3	2.44	0.53
23:BB:483:A:H5''	46:BU:46:LYS:CG	2.39	0.53
30:BY:6:ILE:O	30:BY:34:THR:HA	2.09	0.53
1:CA:1417:G:H1'	1:CA:1483:A:N6	2.24	0.53
1:CA:1453:G:H2'	1:CA:1454:G:O4'	2.09	0.53
1:CA:236:A:H2'	1:CA:237:G:H8	1.74	0.53
1:CA:634:C:H2'	1:CA:635:A:H8	1.74	0.53
1:CA:842:U:H2'	1:CA:843:U:O3'	2.09	0.53
3:CD:148:ALA:C	3:CD:150:LYS:H	2.10	0.53
4:CE:45:VAL:HG11	4:CE:117:ALA:HB2	1.90	0.53
6:CG:75:LYS:HA	6:CG:75:LYS:HZ3	1.73	0.53
9:CJ:37:ARG:CZ	9:CJ:37:ARG:HA	2.39	0.53
10:CK:75:GLU:N	10:CK:75:GLU:CD	2.62	0.53
10:CK:88:PRO:HD3	19:CU:28:LEU:CD1	2.39	0.53
12:CM:108:ARG:HG3	12:CM:108:ARG:HH11	1.73	0.53
23:DB:1241:A:H2'	23:DB:1242:U:H5'	1.90	0.53
23:DB:1258:U:H2'	23:DB:1259:G:C8	2.44	0.53
23:DB:1819:A:OP1	25:DC:154:ALA:HA	2.08	0.53
23:DB:1994:C:O2'	23:DB:1995:U:H5'	2.08	0.53
23:DB:2331:G:H5'	52:DW:39:GLN:HB2	1.91	0.53
23:DB:2352:A:N1	52:DW:30:VAL:HG11	2.24	0.53
23:DB:279:A:H2'	23:DB:280:U:O4'	2.09	0.53
23:DB:2881:U:O2'	23:DB:2882:A:H5'	2.09	0.53
25:DC:90:ILE:HA	25:DC:103:ILE:O	2.09	0.53
42:DN:49:GLU:HB2	42:DN:50:PRO:HD3	1.89	0.53
28:DP:6:GLN:NE2	28:DP:7:LEU:HD12	2.24	0.53
49:DR:38:VAL:HG13	49:DR:54:VAL:HG12	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1487:G:O2'	1:AA:1488:G:H5'	2.09	0.53
18:AB:37:VAL:HG22	18:AB:38:HIS:O	2.09	0.53
2:AC:167:TYR:O	2:AC:168:ARG:HB2	2.08	0.53
11:AL:106:VAL:HA	11:AL:107:LYS:NZ	2.23	0.53
11:AL:79:ILE:HD12	11:AL:96:THR:CG2	2.38	0.53
16:AS:34:SER:O	16:AS:37:SER:HB2	2.09	0.53
23:BB:1018:U:O2'	23:BB:1019:U:H5'	2.09	0.53
23:BB:1229:C:H2'	23:BB:1230:A:C8	2.43	0.53
23:BB:1647:U:H3'	23:BB:1647:U:P	2.49	0.53
23:BB:1672:A:H2'	23:BB:1673:G:C8	2.43	0.53
23:BB:1965:C:H5'	23:BB:1966:A:H2'	1.90	0.53
23:BB:2023:C:O2'	23:BB:2024:G:H5'	2.09	0.53
23:BB:315:G:H2'	23:BB:316:C:C6	2.44	0.53
23:BB:37:C:O2'	23:BB:38:A:H5'	2.08	0.53
23:BB:523:C:H4'	23:BB:540:C:O2	2.09	0.53
23:BB:813:U:H2'	23:BB:814:C:C6	2.44	0.53
25:BC:245:THR:O	25:BC:247:TRP:N	2.41	0.53
47:BF:87:LYS:HG3	47:BF:88:VAL:H	1.73	0.53
48:BG:17:LYS:O	48:BG:23:ILE:HA	2.08	0.53
40:BH:73:ASN:HA	40:BH:141:LYS:O	2.08	0.53
27:BK:19:VAL:CG1	27:BK:43:ILE:HA	2.37	0.53
38:BM:17:ASN:HD21	38:BM:95:LEU:CD1	2.21	0.53
38:BM:71:LYS:HG2	38:BM:93:VAL:HG12	1.90	0.53
42:BN:103:ARG:HB2	42:BN:110:MET:HE2	1.91	0.53
42:BN:12:ARG:HG3	42:BN:13:ASN:H	1.73	0.53
28:BP:77:SER:OG	28:BP:79:VAL:HG22	2.09	0.53
46:BU:26:ASN:N	46:BU:26:ASN:ND2	2.56	0.53
30:BY:47:ILE:HD12	30:BY:54:VAL:HG21	1.90	0.53
1:CA:1227:A:H2'	1:CA:1228:C:O4'	2.09	0.53
1:CA:1480:A:H2'	1:CA:1481:U:C6	2.44	0.53
1:CA:162:A:H2'	1:CA:163:C:O4'	2.09	0.53
1:CA:542:G:O2'	1:CA:543:U:H5'	2.09	0.53
8:CI:11:ARG:NE	8:CI:12:LYS:HB2	2.24	0.53
11:CL:79:ILE:HD12	11:CL:96:THR:CG2	2.38	0.53
32:D4:36:ARG:HG2	32:D4:37:GLN:N	2.22	0.53
23:DB:1425:G:H2'	23:DB:1426:G:C8	2.44	0.53
23:DB:1729:U:H5	23:DB:1731:G:N2	2.07	0.53
23:DB:1999:C:O2'	23:DB:2000:C:H5'	2.09	0.53
23:DB:2537:U:H2'	23:DB:2538:C:H6	1.71	0.53
23:DB:596:U:H2'	23:DB:597:G:H8	1.73	0.53
23:DB:644:A:H2'	23:DB:644:A:N3	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:753:A:H2'	23:DB:754:U:H6	1.72	0.53
23:DB:969:G:OP1	30:DY:17:PRO:HG3	2.09	0.53
29:DE:58:LYS:CD	29:DE:58:LYS:N	2.71	0.53
47:DF:155:ILE:HG22	47:DF:157:THR:N	2.22	0.53
40:DH:14:SER:HB2	40:DH:17:ASP:CB	2.38	0.53
38:DM:17:ASN:HD21	38:DM:95:LEU:HD11	1.74	0.53
1:AA:1089:G:C2'	1:AA:1090:U:H5'	2.37	0.53
1:AA:1110:A:N6	1:AA:1111:A:N6	2.57	0.53
1:AA:124:C:O2'	1:AA:125:U:H5'	2.08	0.53
1:AA:1270:G:H2'	1:AA:1271:A:H8	1.74	0.53
1:AA:229:U:H2'	1:AA:230:G:C8	2.44	0.53
3:AD:60:VAL:HA	3:AD:63:ILE:HD12	1.90	0.53
4:AE:76:ASN:HB2	4:AE:81:GLN:NE2	2.24	0.53
5:AF:47:LEU:HG	5:AF:56:LYS:HA	1.90	0.53
9:AJ:56:HIS:O	9:AJ:57:VAL:HG12	2.09	0.53
23:BB:1468:U:HO2'	23:BB:1469:A:H8	1.55	0.53
23:BB:198:C:O5'	23:BB:198:C:H6	1.92	0.53
23:BB:2458:G:H2'	23:BB:2458:G:N3	2.22	0.53
23:BB:2529:G:H4'	48:BG:174:LYS:CG	2.38	0.53
23:BB:2553:G:H2'	23:BB:2554:U:O4'	2.09	0.53
23:BB:2630:G:O2'	23:BB:2631:G:H5'	2.08	0.53
23:BB:483:A:H5'	46:BU:44:HIS:O	2.09	0.53
23:BB:496:G:C1'	45:BS:61:ASN:HD21	2.22	0.53
23:BB:4:U:H2'	23:BB:5:A:C8	2.44	0.53
23:BB:549:G:H3'	23:BB:549:G:OP2	2.09	0.53
40:BH:4:ILE:HG23	40:BH:16:GLY:HA2	1.90	0.53
40:BH:4:ILE:O	40:BH:6:LEU:HD12	2.09	0.53
38:BM:111:GLU:HA	38:BM:114:ARG:HH22	1.74	0.53
38:BM:40:ARG:HB2	38:BM:93:VAL:CG2	2.39	0.53
45:BS:73:LYS:HE3	45:BS:74:ILE:H	1.74	0.53
39:BX:39:GLN:HB3	39:BX:42:LEU:HD13	1.90	0.53
1:CA:1439:G:H2'	1:CA:1440:U:O4'	2.08	0.53
18:CB:125:PHE:CZ	18:CB:137:THR:HG23	2.44	0.53
2:CC:178:ARG:O	2:CC:178:ARG:HG2	2.08	0.53
7:CH:50:VAL:HG22	7:CH:50:VAL:O	2.09	0.53
1:CA:878:A:H5''	7:CH:80:PRO:HG2	1.89	0.53
10:CK:53:GLY:O	10:CK:56:LYS:HB3	2.09	0.53
11:CL:81:ILE:HA	11:CL:95:HIS:O	2.09	0.53
21:CN:50:LEU:CD2	21:CN:51:PRO:HD3	2.38	0.53
16:CS:7:GLY:H	16:CS:8:PRO:HD3	1.73	0.53
22:DA:35:C:H2'	22:DA:36:C:O4'	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1313:U:O2	23:DB:1313:U:H2'	2.08	0.53
23:DB:1346:G:O2'	23:DB:1347:A:H5'	2.09	0.53
23:DB:1430:G:H2'	23:DB:1431:A:C8	2.44	0.53
23:DB:1734:G:O2'	23:DB:1735:A:H5'	2.09	0.53
23:DB:2106:U:H2'	23:DB:2107:G:O4'	2.09	0.53
23:DB:2564:A:OP1	23:DB:2648:G:H4'	2.09	0.53
23:DB:2893:A:H4'	23:DB:2894:G:O5'	2.08	0.53
25:DC:144:GLU:OE2	25:DC:188:ARG:HG3	2.09	0.53
25:DC:262:THR:O	25:DC:263:ASP:C	2.47	0.53
26:DD:119:ALA:HB1	26:DD:163:GLY:N	2.22	0.53
26:DD:56:LYS:HG3	26:DD:58:ASN:HB3	1.91	0.53
47:DF:29:ARG:HE	47:DF:29:ARG:H	1.57	0.53
47:DF:35:LEU:HD23	47:DF:36:ASN:H	1.73	0.53
48:DG:24:THR:HG22	48:DG:34:ARG:HA	1.90	0.53
42:DN:22:ARG:CZ	42:DN:70:THR:H	2.21	0.53
52:DW:9:THR:C	52:DW:10:ARG:HD3	2.29	0.53
52:DW:38:ARG:N	52:DW:56:HIS:HD2	2.07	0.53
30:DY:8:GLN:HB3	30:DY:31:ILE:O	2.08	0.53
1:AA:201:G:H2'	1:AA:202:G:O4'	2.09	0.53
1:AA:778:G:H2'	1:AA:779:C:C6	2.44	0.53
2:AC:149:LYS:HE3	2:AC:166:TRP:CH2	2.44	0.53
9:AJ:44:THR:HG23	9:AJ:69:THR:O	2.09	0.53
9:AJ:6:ILE:O	9:AJ:8:ILE:HD12	2.09	0.53
11:AL:110:LYS:O	11:AL:113:ARG:HG3	2.09	0.53
12:AM:21:ILE:HB	12:AM:24:VAL:CG2	2.37	0.53
17:AT:53:MET:HA	17:AT:56:ILE:CD1	2.39	0.53
23:BB:1060:U:O2	23:BB:1088:A:C8	2.62	0.53
23:BB:1640:A:H2'	23:BB:1641:A:H8	1.74	0.53
23:BB:18:U:H2'	23:BB:19:A:C8	2.44	0.53
23:BB:2455:G:H2'	23:BB:2456:C:H6	1.74	0.53
23:BB:2645:G:C5'	23:BB:2732:G:H8	2.21	0.53
23:BB:734:A:H2'	23:BB:735:A:O4'	2.08	0.53
23:BB:742:A:H2'	23:BB:743:A:C8	2.44	0.53
23:BB:873:C:H4'	38:BM:64:TRP:NE1	2.22	0.53
47:BF:62:GLN:NE2	47:BF:91:ARG:HE	2.07	0.53
41:BJ:59:ALA:HB1	41:BJ:101:ILE:HD11	1.90	0.53
41:BJ:54:ILE:HD12	41:BJ:55:ILE:N	2.23	0.53
37:BL:3:LEU:O	37:BL:5:THR:HG23	2.08	0.53
35:BV:38:LEU:HG	35:BV:40:ILE:HG23	1.90	0.53
1:CA:1481:U:O2'	1:CA:1482:G:H5'	2.08	0.53
1:CA:6:G:H4'	1:CA:298:A:H4'	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:802:A:H2'	1:CA:803:G:O4'	2.08	0.53
18:CB:209:VAL:HG23	18:CB:210:THR:N	2.23	0.53
18:CB:20:ARG:HA	18:CB:38:HIS:CE1	2.44	0.53
18:CB:216:VAL:HG23	18:CB:217:ALA:H	1.74	0.53
2:CC:71:ARG:O	2:CC:74:ILE:HG22	2.09	0.53
3:CD:59:LYS:O	3:CD:63:ILE:HG13	2.09	0.53
5:CF:68:GLN:HG2	5:CF:69:GLU:N	2.24	0.53
17:CT:67:HIS:CG	17:CT:68:LYS:H	2.27	0.53
23:DB:1021:A:H61	23:DB:1142:A:H61	1.57	0.53
23:DB:1210:G:H5'	23:DB:1212:G:O4'	2.09	0.53
23:DB:1431:A:H2'	23:DB:1432:G:C8	2.43	0.53
23:DB:1872:A:H8	23:DB:1872:A:O5'	1.91	0.53
23:DB:1873:G:O2'	23:DB:1874:C:H5'	2.09	0.53
23:DB:579:G:H4'	23:DB:2018:G:H5''	1.90	0.53
23:DB:2296:U:H4'	23:DB:2297:A:OP1	2.09	0.53
23:DB:2405:G:H1'	23:DB:2412:A:H61	1.73	0.53
23:DB:2651:C:O2'	23:DB:2652:C:H5'	2.09	0.53
23:DB:2875:C:H2'	23:DB:2876:G:C8	2.43	0.53
23:DB:453:A:H5''	56:DB:3501:HOH:O	2.08	0.53
23:DB:697:G:H2'	23:DB:698:C:C6	2.44	0.53
25:DC:77:VAL:HG22	25:DC:113:ASP:O	2.09	0.53
29:DE:48:THR:O	29:DE:52:VAL:HG23	2.07	0.53
47:DF:100:GLU:OE1	47:DF:100:GLU:HA	2.09	0.53
27:DK:109:SER:OG	27:DK:111:LYS:HG2	2.09	0.53
27:DK:87:LEU:HB2	27:DK:93:GLN:C	2.29	0.53
38:DM:40:ARG:HB2	38:DM:93:VAL:CG2	2.39	0.53
42:DN:45:ARG:O	42:DN:49:GLU:HG3	2.09	0.53
49:DR:98:ILE:O	49:DR:99:THR:HG23	2.09	0.53
46:DU:64:ILE:HG13	46:DU:65:GLN:N	2.23	0.53
1:AA:1085:U:H3'	1:AA:1086:U:C5	2.44	0.52
1:AA:1469:C:H2'	1:AA:1470:U:O4'	2.08	0.52
1:AA:50:A:N6	1:AA:361:G:H4'	2.24	0.52
1:AA:699:C:C2'	1:AA:700:G:H5''	2.39	0.52
1:AA:842:U:OP1	1:AA:843:U:H4'	2.09	0.52
18:AB:209:VAL:HG23	18:AB:210:THR:N	2.10	0.52
2:AC:102:ILE:H	2:AC:102:ILE:HD12	1.73	0.52
2:AC:115:VAL:O	2:AC:119:ILE:HG22	2.09	0.52
2:AC:13:ILE:C	2:AC:15:LYS:H	2.12	0.52
4:AE:14:LEU:HD22	4:AE:15:ILE:N	2.24	0.52
7:AH:108:GLY:O	7:AH:110:MET:HG3	2.09	0.52
10:AK:95:THR:HG23	10:AK:96:ILE:N	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AO:77:ARG:HH11	20:AO:77:ARG:HB3	1.74	0.52
32:B4:33:HIS:O	32:B4:35:GLN:HG3	2.09	0.52
23:BB:1199:U:H2'	23:BB:1200:C:H6	1.73	0.52
23:BB:1260:A:O2'	23:BB:1261:C:H5'	2.08	0.52
23:BB:1930:G:H2'	23:BB:1968:G:C6	2.44	0.52
23:BB:2290:G:H2'	23:BB:2291:U:C6	2.44	0.52
23:BB:2300:C:H2'	23:BB:2301:C:C6	2.43	0.52
23:BB:30:G:H2'	23:BB:31:C:H6	1.70	0.52
23:BB:545:U:H1'	23:BB:548:G:O6	2.09	0.52
23:BB:571:U:N3	23:BB:575:A:N7	2.57	0.52
23:BB:584:C:OP2	44:BQ:5:ARG:HD3	2.09	0.52
23:BB:750:A:H2'	23:BB:751:A:H5''	1.91	0.52
26:BD:111:GLY:H	26:BD:194:PRO:CG	2.23	0.52
47:BF:45:ASP:O	47:BF:47:LYS:N	2.41	0.52
47:BF:45:ASP:O	47:BF:46:LYS:HE2	2.10	0.52
27:BK:34:GLY:O	27:BK:36:GLY:N	2.42	0.52
46:BU:33:VAL:HG13	46:BU:66:VAL:HB	1.91	0.52
46:BU:12:VAL:HG22	46:BU:69:VAL:HG12	1.91	0.52
1:CA:384:G:H2'	1:CA:385:C:H6	1.74	0.52
1:CA:562:U:H4'	1:CA:563:A:O5'	2.09	0.52
18:CB:172:ILE:HG22	18:CB:176:ASN:HD21	1.73	0.52
18:CB:86:CYS:N	18:CB:88:GLN:HE21	2.07	0.52
5:CF:47:LEU:HG	5:CF:56:LYS:HA	1.90	0.52
7:CH:37:ASN:O	7:CH:40:LYS:HB3	2.09	0.52
12:CM:10:ASP:HA	12:CM:44:ILE:HD13	1.90	0.52
20:CO:71:LYS:HZ3	20:CO:71:LYS:HB3	1.74	0.52
22:DA:28:C:H2'	22:DA:29:A:O4'	2.08	0.52
22:DA:87:U:H2'	22:DA:88:C:C5'	2.39	0.52
23:DB:1623:G:O2'	23:DB:1624:U:H5'	2.09	0.52
23:DB:2530:A:H2'	23:DB:2531:A:H5''	1.91	0.52
23:DB:2605:U:H2'	23:DB:2606:C:C6	2.43	0.52
23:DB:2615:U:C2	31:D0:3:GLN:HA	2.44	0.52
23:DB:372:G:OP2	51:DZ:61:LYS:HE2	2.09	0.52
23:DB:39:G:H2'	23:DB:40:U:C6	2.44	0.52
23:DB:99:U:O2	23:DB:99:U:O4'	2.27	0.52
47:DF:116:LEU:HG	47:DF:117:SER:H	1.73	0.52
24:DI:54:ILE:HD13	24:DI:55:PRO:N	2.23	0.52
50:DT:39:THR:CG2	50:DT:42:GLU:H	2.23	0.52
50:DT:44:LYS:O	50:DT:48:GLN:HG2	2.10	0.52
50:DT:50:LEU:HD22	50:DT:50:LEU:N	2.23	0.52
46:DU:13:LEU:HA	46:DU:18:LYS:CE	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:DZ:5:CYS:HB3	51:DZ:9:GLY:H	1.73	0.52
1:AA:1117:A:H4'	8:AI:105:ARG:NH1	2.24	0.52
1:AA:1125:U:H2'	1:AA:1126:U:C5'	2.36	0.52
1:AA:218:U:H2'	1:AA:219:U:H6	1.74	0.52
1:AA:601:G:H2'	1:AA:602:A:H8	1.74	0.52
18:AB:131:LYS:HZ3	18:AB:132:GLU:N	2.05	0.52
3:AD:54:LEU:O	3:AD:54:LEU:HD22	2.09	0.52
5:AF:18:VAL:O	5:AF:22:ILE:HG13	2.09	0.52
6:AG:28:ILE:HG21	6:AG:101:ARG:HG2	1.90	0.52
1:AA:1148:U:H5'	8:AI:6:TYR:OH	2.08	0.52
12:AM:106:ARG:NE	12:AM:112:ARG:HG2	2.24	0.52
20:AO:36:ILE:C	20:AO:38:HIS:H	2.12	0.52
17:AT:43:LYS:HB3	17:AT:86:ALA:CB	2.39	0.52
34:B3:14:LYS:O	34:B3:21:PHE:HA	2.10	0.52
22:BA:50:A:OP1	43:BO:68:LYS:HB2	2.09	0.52
23:BB:1241:A:H2'	23:BB:1242:U:H5'	1.91	0.52
23:BB:1904:G:O2'	23:BB:1905:C:H5'	2.09	0.52
23:BB:4:U:H2'	23:BB:5:A:H8	1.74	0.52
23:BB:725:G:H2'	23:BB:726:G:O4'	2.09	0.52
29:BE:105:LEU:HD21	29:BE:177:PRO:HA	1.90	0.52
47:BF:133:GLU:HA	47:BF:150:GLY:CA	2.39	0.52
24:BI:2:LYS:NZ	24:BI:2:LYS:HB3	2.24	0.52
27:BK:87:LEU:HB2	27:BK:93:GLN:C	2.29	0.52
23:BB:2485:G:H5''	38:BM:125:PRO:HG3	1.91	0.52
42:BN:38:LEU:HD11	42:BN:42:LYS:HD2	1.91	0.52
42:BN:45:ARG:O	42:BN:49:GLU:HG3	2.09	0.52
45:BS:68:ASP:C	45:BS:69:LEU:HD12	2.30	0.52
50:BT:8:LEU:C	50:BT:9:LYS:HE2	2.29	0.52
46:BU:43:LYS:HD3	46:BU:44:HIS:N	2.24	0.52
1:CA:1225:A:O2'	16:CS:77:ARG:HD3	2.08	0.52
1:CA:1240:U:O2'	6:CG:31:VAL:HG11	2.08	0.52
1:CA:1332:A:H2'	1:CA:1333:A:C8	2.44	0.52
1:CA:135:C:O2	13:CP:1:MET:HB2	2.10	0.52
1:CA:539:A:H2'	1:CA:540:G:H8	1.75	0.52
1:CA:570:G:H2'	1:CA:571:U:C6	2.45	0.52
6:CG:90:VAL:CG2	6:CG:95:ARG:HG3	2.40	0.52
8:CI:11:ARG:HD3	8:CI:73:GLY:HA2	1.89	0.52
1:CA:1328:C:OP1	12:CM:27:THR:HG21	2.09	0.52
12:CM:52:ILE:HG23	12:CM:53:ASP:N	2.23	0.52
21:CN:1:ALA:HB1	21:CN:6:LYS:HE3	1.91	0.52
16:CS:25:GLY:O	16:CS:27:LYS:HE2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1320:C:C2	16:CS:71:GLY:HA3	2.45	0.52
33:D1:5:ARG:NH1	33:D1:25:ASN:HB2	2.23	0.52
23:DB:1013:C:H2'	23:DB:1014:A:C8	2.43	0.52
23:DB:1099:G:C8	24:DI:3:LYS:HB2	2.44	0.52
23:DB:1210:G:N3	23:DB:1212:G:N2	2.57	0.52
23:DB:1281:G:H2'	23:DB:1282:U:C6	2.44	0.52
23:DB:1488:C:O2'	23:DB:1489:C:H5'	2.10	0.52
23:DB:1771:C:H2'	23:DB:1772:A:C8	2.44	0.52
23:DB:2104:C:C6	23:DB:2104:C:C3'	2.92	0.52
23:DB:2300:C:H2'	23:DB:2301:C:C6	2.44	0.52
23:DB:2471:A:O2'	23:DB:2472:G:C8	2.58	0.52
23:DB:850:U:H2'	23:DB:851:C:H6	1.73	0.52
47:DF:12:VAL:O	47:DF:16:MET:HG2	2.09	0.52
48:DG:17:LYS:O	48:DG:23:ILE:HA	2.09	0.52
27:DK:64:ARG:HD2	27:DK:102:PRO:O	2.09	0.52
37:DL:51:GLU:OE1	37:DL:57:LEU:HB2	2.10	0.52
37:DL:65:GLY:O	37:DL:66:PHE:HB3	2.09	0.52
38:DM:42:THR:H	38:DM:45:GLN:CD	2.12	0.52
28:DP:19:PHE:CE2	28:DP:25:VAL:HG11	2.44	0.52
50:DT:15:HIS:N	50:DT:32:LEU:HA	2.15	0.52
52:DW:37:VAL:CG1	52:DW:38:ARG:H	2.15	0.52
1:AA:191:G:H8	1:AA:191:G:OP2	1.92	0.52
1:AA:143:A:H2	1:AA:220:G:H22	1.56	0.52
1:AA:648:A:H2'	1:AA:649:A:H8	1.75	0.52
18:AB:102:ASN:O	18:AB:106:VAL:HG23	2.09	0.52
18:AB:134:LEU:HA	18:AB:137:THR:HG1	1.75	0.52
3:AD:148:ALA:C	3:AD:150:LYS:H	2.12	0.52
7:AH:124:ILE:HG22	7:AH:125:ILE:N	2.24	0.52
8:AI:56:MET:C	8:AI:58:GLU:H	2.12	0.52
14:AQ:18:LYS:HE3	14:AQ:48:GLU:HG2	1.89	0.52
10:AK:88:PRO:HD3	19:AU:28:LEU:CD1	2.39	0.52
19:AU:34:ARG:HD3	19:AU:39:LYS:CE	2.35	0.52
19:AU:40:PRO:C	19:AU:42:THR:H	2.13	0.52
34:B3:60:CYS:O	34:B3:61:LEU:HD23	2.08	0.52
32:B4:24:ARG:NH2	32:B4:36:ARG:HG3	2.23	0.52
22:BA:64:G:H2'	22:BA:65:U:C6	2.44	0.52
23:BB:1021:A:H61	23:BB:1142:A:H61	1.56	0.52
23:BB:1199:U:H1'	44:BQ:2:ARG:O	2.09	0.52
23:BB:2671:G:H2'	23:BB:2672:U:H6	1.74	0.52
23:BB:2677:G:H2'	23:BB:2678:C:H6	1.74	0.52
23:BB:2769:U:H2'	23:BB:2770:G:C8	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:346:A:H5'	23:BB:346:A:N3	2.24	0.52
23:BB:643:A:C5	23:BB:644:A:N7	2.78	0.52
23:BB:851:C:H2'	23:BB:852:U:C6	2.45	0.52
23:BB:85:G:N2	23:BB:98:G:H1'	2.24	0.52
25:BC:12:ARG:HA	25:BC:15:VAL:HG23	1.91	0.52
25:BC:188:ARG:HH21	25:BC:188:ARG:HG2	1.74	0.52
25:BC:209:ALA:HA	25:BC:212:TRP:CE2	2.45	0.52
29:BE:116:ASP:O	29:BE:185:LYS:HE3	2.08	0.52
47:BF:62:GLN:NE2	47:BF:90:LEU:HA	2.24	0.52
40:BH:116:ARG:CZ	40:BH:131:SER:HB2	2.39	0.52
27:BK:75:SER:HA	28:BP:72:VAL:O	2.09	0.52
38:BM:102:LEU:H	38:BM:102:LEU:HD22	1.72	0.52
35:BV:75:GLN:HB3	35:BV:90:ASP:HB2	1.91	0.52
52:BW:46:ALA:HB2	52:BW:78:PHE:CD1	2.45	0.52
1:CA:1320:C:OP2	16:CS:69:LYS:HE3	2.09	0.52
2:CC:81:GLU:O	2:CC:85:LYS:HB2	2.09	0.52
10:CK:51:PHE:HZ	10:CK:61:ALA:HA	1.74	0.52
12:CM:55:LEU:O	12:CM:59:VAL:HG12	2.09	0.52
14:CQ:76:ARG:HH12	14:CQ:78:VAL:HA	1.74	0.52
23:DB:137:U:H2'	23:DB:138:U:C1'	2.40	0.52
23:DB:1439:A:N7	23:DB:1440:U:N1	2.57	0.52
23:DB:1777:U:O2'	23:DB:1778:U:H5'	2.09	0.52
23:DB:2179:C:H2'	23:DB:2179:C:O2	2.08	0.52
23:DB:2769:U:H2'	23:DB:2770:G:C8	2.45	0.52
23:DB:523:C:H4'	23:DB:540:C:O2	2.09	0.52
23:DB:813:U:H2'	23:DB:814:C:C6	2.44	0.52
23:DB:871:U:H1'	23:DB:907:G:N1	2.25	0.52
23:DB:936:A:H2'	23:DB:937:C:C6	2.45	0.52
25:DC:132:ARG:NH2	25:DC:169:ALA:HA	2.24	0.52
23:DB:1812:U:H1'	25:DC:43:ASN:HD21	1.74	0.52
48:DG:89:VAL:HG23	48:DG:160:GLY:O	2.10	0.52
23:DB:1098:A:C3'	24:DI:4:VAL:N	2.72	0.52
41:DJ:104:ALA:O	41:DJ:108:MET:HG3	2.10	0.52
41:DJ:44:TYR:CE2	44:DQ:59:LEU:HD11	2.45	0.52
27:DK:87:LEU:HD12	27:DK:93:GLN:H	1.73	0.52
23:DB:825:A:O2'	37:DL:54:GLN:HB3	2.09	0.52
30:DY:3:THR:HB	30:DY:36:GLU:CG	2.39	0.52
1:AA:1451:U:H5''	1:AA:1452:C:H5	1.74	0.52
1:AA:295:C:H2'	1:AA:296:U:C6	2.45	0.52
1:AA:555:U:H2'	1:AA:556:C:C6	2.43	0.52
1:AA:6:G:H4'	1:AA:298:A:H4'	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:882:C:O2'	1:AA:883:C:H5'	2.09	0.52
18:AB:93:HIS:O	18:AB:94:ARG:HG2	2.10	0.52
4:AE:85:LYS:HG3	4:AE:93:VAL:O	2.09	0.52
7:AH:37:ASN:O	7:AH:41:GLU:HG2	2.09	0.52
8:AI:18:VAL:HG21	8:AI:82:ILE:N	2.23	0.52
8:AI:80:HIS:O	8:AI:84:ARG:HG3	2.09	0.52
9:AJ:53:ILE:CG2	9:AJ:61:ALA:HB1	2.31	0.52
20:AO:3:LEU:HG	20:AO:8:THR:OG1	2.10	0.52
13:AP:75:ILE:HG22	13:AP:80:LYS:HD2	1.90	0.52
23:BB:1097:U:C2'	23:BB:1098:A:H5'	2.38	0.52
23:BB:1244:A:O2'	23:BB:1245:G:H5'	2.09	0.52
23:BB:1545:A:H2'	23:BB:1546:G:O4'	2.08	0.52
23:BB:1838:C:H4'	23:BB:1839:G:H8	1.73	0.52
23:BB:1997:C:H2'	23:BB:1998:A:C8	2.44	0.52
23:BB:2443:C:O2'	23:BB:2444:G:H5'	2.10	0.52
23:BB:878:A:H3'	23:BB:878:A:OP1	2.09	0.52
23:BB:90:U:H3'	23:BB:91:A:C5'	2.38	0.52
25:BC:262:THR:O	25:BC:263:ASP:C	2.48	0.52
25:BC:52:HIS:N	25:BC:52:HIS:CD2	2.76	0.52
47:BF:41:GLU:HG3	47:BF:48:LEU:HD11	1.91	0.52
48:BG:118:ALA:C	48:BG:120:ILE:H	2.13	0.52
48:BG:34:ARG:HD3	48:BG:34:ARG:N	2.24	0.52
40:BH:40:THR:O	40:BH:42:LYS:N	2.39	0.52
40:BH:58:LEU:HA	40:BH:61:VAL:HG12	1.91	0.52
38:BM:21:ALA:HB2	38:BM:100:LYS:HG2	1.92	0.52
30:BY:3:THR:HB	30:BY:36:GLU:CG	2.40	0.52
30:BY:6:ILE:HD13	30:BY:6:ILE:N	2.22	0.52
1:CA:1115:U:H2'	1:CA:1116:U:C6	2.44	0.52
1:CA:1527:U:O2'	1:CA:1528:U:H5'	2.09	0.52
1:CA:1503:A:C8	1:CA:1531:A:H1'	2.44	0.52
1:CA:179:A:H2'	1:CA:180:U:O4'	2.08	0.52
1:CA:476:U:H2'	1:CA:477:C:C6	2.43	0.52
1:CA:642:A:H2'	1:CA:643:C:C6	2.44	0.52
1:CA:749:A:O2'	1:CA:750:C:H5'	2.09	0.52
18:CB:125:PHE:HZ	18:CB:137:THR:HG23	1.74	0.52
7:CH:55:LYS:HE3	7:CH:56:PRO:HD2	1.91	0.52
12:CM:38:ILE:CG2	12:CM:42:VAL:HG21	2.40	0.52
12:CM:69:ARG:O	12:CM:70:ARG:C	2.48	0.52
14:CQ:60:ILE:HG22	14:CQ:74:LEU:HA	1.91	0.52
17:CT:73:ARG:HG3	17:CT:74:HIS:N	2.24	0.52
36:D2:3:ARG:CZ	36:D2:3:ARG:HA	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:D4:7:VAL:HG13	32:D4:8:LYS:N	2.14	0.52
23:DB:1064:C:H5'	24:DI:88:GLY:HA3	1.90	0.52
23:DB:1782:U:H2'	23:DB:1783:A:H5'	1.90	0.52
23:DB:1916:A:H2'	23:DB:1917:U:O4'	2.08	0.52
23:DB:917:A:H5''	23:DB:2268:A:H61	1.75	0.52
23:DB:1629:U:O2	23:DB:2698:U:H5''	2.09	0.52
23:DB:736:C:H2'	23:DB:737:C:C6	2.44	0.52
23:DB:1190:G:OP1	37:DL:32:GLY:HA2	2.08	0.52
38:DM:21:ALA:HB2	38:DM:100:LYS:HG2	1.91	0.52
52:DW:49:ASN:C	52:DW:50:VAL:HG22	2.30	0.52
39:DX:41:HIS:O	39:DX:44:LYS:HB3	2.09	0.52
51:DZ:32:ASN:C	51:DZ:33:LEU:HD12	2.30	0.52
1:AA:1361:G:C2'	1:AA:1362:A:H5''	2.29	0.52
1:AA:137:U:H2'	1:AA:138:G:H8	1.74	0.52
1:AA:370:C:H2'	1:AA:371:A:H8	1.73	0.52
1:AA:461:A:H3'	1:AA:462:G:O4'	2.09	0.52
1:AA:577:G:O2'	1:AA:578:C:H5'	2.09	0.52
1:AA:755:G:O2'	1:AA:756:C:H5'	2.10	0.52
18:AB:38:HIS:O	18:AB:39:ILE:HD13	2.09	0.52
4:AE:96:GLN:HB3	4:AE:123:LEU:HD12	1.90	0.52
4:AE:98:ALA:HB2	4:AE:123:LEU:HG	1.90	0.52
6:AG:114:SER:O	6:AG:118:ARG:HG3	2.08	0.52
9:AJ:28:THR:HG21	9:AJ:90:LEU:HD23	1.90	0.52
21:AN:17:ASP:O	21:AN:21:ALA:HB3	2.09	0.52
34:B3:31:ILE:CD1	34:B3:34:LYS:HD3	2.39	0.52
22:BA:28:C:H2'	22:BA:29:A:O4'	2.10	0.52
23:BB:1439:A:H1'	23:BB:1553:A:N6	2.25	0.52
23:BB:1661:G:O2'	23:BB:1662:U:H5'	2.10	0.52
23:BB:1946:U:H2'	23:BB:1947:C:H6	1.74	0.52
23:BB:2052:A:OP1	26:BD:145:SER:HA	2.09	0.52
23:BB:2199:A:H5'	23:BB:2200:C:OP2	2.10	0.52
23:BB:2753:A:H2	32:B4:15:LYS:HZ1	1.57	0.52
23:BB:2885:G:H2'	23:BB:2886:A:O4'	2.09	0.52
23:BB:692:C:H2'	23:BB:693:A:C8	2.45	0.52
23:BB:833:A:H2'	23:BB:834:G:C8	2.45	0.52
23:BB:863:A:H2'	23:BB:864:G:H8	1.73	0.52
23:BB:1820:U:OP1	25:BC:176:ARG:HD2	2.09	0.52
25:BC:28:PRO:HG2	25:BC:33:LEU:HD11	1.92	0.52
26:BD:106:LYS:N	26:BD:106:LYS:HD3	2.25	0.52
47:BF:32:LYS:H	47:BF:95:MET:HE1	1.74	0.52
24:BI:48:ILE:HG22	24:BI:49:GLU:HG2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BP:91:VAL:O	28:BP:92:ARG:HB3	2.10	0.52
35:BV:10:LYS:HG2	35:BV:11:GLU:HG3	1.90	0.52
52:BW:35:ILE:HG13	52:BW:57:THR:OG1	2.08	0.52
1:CA:1412:C:H2'	1:CA:1413:A:C8	2.45	0.52
18:CB:52:ALA:HB2	18:CB:197:PHE:HB3	1.92	0.52
18:CB:86:CYS:O	18:CB:88:GLN:N	2.41	0.52
2:CC:120:THR:HG23	2:CC:197:VAL:CG2	2.40	0.52
3:CD:145:ARG:HB3	3:CD:147:LYS:HG3	1.92	0.52
6:CG:112:ASP:H	6:CG:118:ARG:CZ	2.22	0.52
4:CE:156:ARG:HB2	7:CH:43:GLY:HA3	1.91	0.52
11:CL:5:GLN:HA	11:CL:8:ARG:NH2	2.24	0.52
12:CM:14:ALA:HB2	12:CM:42:VAL:CG2	2.38	0.52
12:CM:21:ILE:CB	12:CM:24:VAL:HG22	2.31	0.52
12:CM:3:ILE:HG22	12:CM:56:ARG:HA	1.92	0.52
12:CM:82:LEU:HD21	16:CS:64:GLU:OE2	2.09	0.52
13:CP:28:ARG:HD3	13:CP:29:ASN:ND2	2.25	0.52
16:CS:47:THR:HG23	16:CS:60:PHE:CZ	2.44	0.52
23:DB:1100:C:H2'	23:DB:1101:U:C6	2.45	0.52
23:DB:1353:A:H2'	23:DB:1354:A:H8	1.74	0.52
23:DB:1475:G:H1'	23:DB:1476:U:H5	1.75	0.52
23:DB:1582:C:H2'	23:DB:1583:A:O4'	2.09	0.52
23:DB:1593:A:H2'	23:DB:1594:U:O4'	2.09	0.52
23:DB:1676:A:H2'	23:DB:1677:A:O4'	2.10	0.52
23:DB:1711:A:H2'	23:DB:1712:U:C6	2.45	0.52
23:DB:2337:G:H2'	23:DB:2337:G:N3	2.24	0.52
23:DB:2674:G:H2'	23:DB:2675:A:C8	2.44	0.52
23:DB:2884:U:H2'	23:DB:2885:G:C8	2.44	0.52
23:DB:443:A:C4	29:DE:40:ARG:HD3	2.44	0.52
23:DB:734:A:H2'	23:DB:735:A:O4'	2.10	0.52
23:DB:742:A:O2'	23:DB:743:A:H5'	2.10	0.52
26:DD:123:LYS:HD2	26:DD:165:MET:HE1	1.91	0.52
48:DG:104:LEU:HB2	48:DG:112:VAL:HB	1.92	0.52
40:DH:121:VAL:HG12	40:DH:123:ARG:HD2	1.90	0.52
40:DH:125:THR:CA	40:DH:146:VAL:HB	2.31	0.52
41:DJ:19:ASP:HB2	41:DJ:57:LEU:HB2	1.90	0.52
27:DK:118:LEU:O	27:DK:120:PRO:HD2	2.10	0.52
50:DT:8:LEU:C	50:DT:9:LYS:HE2	2.29	0.52
35:DV:10:LYS:HG2	35:DV:11:GLU:HG3	1.91	0.52
1:AA:1264:U:O2'	1:AA:1265:C:H5'	2.09	0.52
1:AA:1492:A:H2'	23:BB:1913:A:C6	2.45	0.52
1:AA:364:A:H2'	1:AA:365:U:O2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:78:A:H2'	1:AA:79:G:C8	2.45	0.52
1:AA:868:C:H2'	1:AA:869:G:O4'	2.09	0.52
1:AA:937:A:H3'	56:AA:1753:HOH:O	2.09	0.52
4:AE:44:ARG:HA	4:AE:71:ILE:O	2.09	0.52
6:AG:135:LYS:C	6:AG:135:LYS:HD3	2.29	0.52
12:AM:18:LEU:HD23	12:AM:18:LEU:O	2.10	0.52
14:AQ:76:ARG:HH12	14:AQ:78:VAL:HA	1.75	0.52
12:AM:92:ARG:HH11	16:AS:79:TYR:HE2	1.58	0.52
17:AT:67:HIS:CG	17:AT:68:LYS:H	2.28	0.52
23:BB:1374:G:H2'	23:BB:1375:U:C6	2.44	0.52
23:BB:2432:A:O2'	23:BB:2433:A:H5'	2.09	0.52
23:BB:2621:G:P	26:BD:124:ARG:HH22	2.32	0.52
23:BB:2834:G:H1'	23:BB:2883:A:N6	2.25	0.52
23:BB:2884:U:H2'	23:BB:2885:G:C8	2.43	0.52
23:BB:531:C:O2'	23:BB:563:A:H5''	2.10	0.52
25:BC:12:ARG:HA	25:BC:15:VAL:CG2	2.40	0.52
29:BE:60:TRP:HE1	29:BE:71:GLY:HA2	1.73	0.52
48:BG:72:ASN:O	48:BG:76:ILE:HG12	2.09	0.52
41:BJ:17:VAL:CG2	41:BJ:137:PRO:HB2	2.34	0.52
41:BJ:44:TYR:CD2	44:BQ:59:LEU:HD21	2.44	0.52
42:BN:9:GLN:O	42:BN:11:ASN:N	2.41	0.52
43:BO:74:VAL:O	43:BO:77:ALA:HB3	2.10	0.52
52:BW:18:LYS:HA	52:BW:36:ILE:HG12	1.90	0.52
51:BZ:33:LEU:H	51:BZ:52:SER:CB	2.23	0.52
1:CA:218:U:H2'	1:CA:219:U:H6	1.75	0.52
1:CA:229:U:H2'	1:CA:230:G:C8	2.44	0.52
1:CA:683:G:O2'	1:CA:684:U:H5'	2.10	0.52
1:CA:735:C:O2'	1:CA:736:C:H5'	2.10	0.52
1:CA:803:G:H2'	1:CA:804:U:H6	1.73	0.52
1:CA:947:G:H2'	1:CA:948:C:C6	2.45	0.52
1:CA:959:A:N6	16:CS:77:ARG:HA	2.24	0.52
5:CF:6:ILE:HG13	5:CF:62:MET:HB2	1.92	0.52
6:CG:25:PHE:HD1	6:CG:100:MET:HG2	1.75	0.52
11:CL:66:ILE:N	11:CL:66:ILE:HD12	2.25	0.52
21:CN:40:ARG:NH2	16:CS:6:LYS:HB2	2.25	0.52
14:CQ:59:GLU:HB3	14:CQ:75:VAL:HG23	1.92	0.52
33:D1:9:LYS:HD3	33:D1:9:LYS:N	2.21	0.52
32:D4:33:HIS:O	32:D4:35:GLN:HG3	2.10	0.52
32:D4:24:ARG:NH2	32:D4:36:ARG:HG3	2.24	0.52
22:DA:113:C:H2'	22:DA:114:C:C6	2.45	0.52
23:DB:1151:A:H2'	23:DB:1152:C:H6	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:125:A:H3'	23:DB:126:A:H5''	1.90	0.52
23:DB:1260:A:O2'	23:DB:1261:C:H5'	2.10	0.52
23:DB:1361:G:H2'	23:DB:1362:C:C6	2.45	0.52
23:DB:1374:G:O2'	23:DB:1375:U:H5'	2.10	0.52
23:DB:2408:U:O2'	23:DB:2409:G:H5'	2.09	0.52
23:DB:2454:G:C2'	23:DB:2455:G:H5'	2.40	0.52
23:DB:627:A:OP1	23:DB:627:A:H2'	2.10	0.52
26:DD:106:LYS:HD3	26:DD:106:LYS:H	1.75	0.52
29:DE:60:TRP:C	29:DE:62:GLN:H	2.12	0.52
47:DF:107:VAL:HG21	47:DF:175:PRO:HG3	1.91	0.52
47:DF:111:ARG:HH11	47:DF:135:ILE:HG21	1.73	0.52
40:DH:40:THR:O	40:DH:42:LYS:N	2.41	0.52
40:DH:4:ILE:O	40:DH:6:LEU:HD12	2.09	0.52
38:DM:114:ARG:HG3	38:DM:130:PHE:CG	2.45	0.52
44:DQ:57:ARG:HA	44:DQ:60:TRP:CE3	2.44	0.52
50:DT:39:THR:O	50:DT:40:LYS:HB3	2.09	0.52
35:DV:53:LYS:HZ3	35:DV:53:LYS:HA	1.75	0.52
30:DY:50:VAL:HB	30:DY:53:MET:CB	2.39	0.52
1:AA:1031:C:H4'	1:AA:1032:G:O5'	2.09	0.52
1:AA:179:A:H2'	1:AA:180:U:O4'	2.09	0.52
1:AA:539:A:H2'	1:AA:540:G:C8	2.44	0.52
1:AA:61:G:H2'	1:AA:62:U:O4'	2.09	0.52
1:AA:821:G:H2'	1:AA:822:U:C6	2.45	0.52
6:AG:12:LEU:HD13	6:AG:13:PRO:HD2	1.92	0.52
11:AL:66:ILE:N	11:AL:66:ILE:HD12	2.25	0.52
12:AM:44:ILE:HD12	12:AM:44:ILE:H	1.75	0.52
1:AA:254:G:OP1	14:AQ:68:LYS:O	2.26	0.52
23:BB:1100:C:H2'	23:BB:1101:U:H6	1.74	0.52
23:BB:1560:G:H2'	23:BB:1561:C:C6	2.43	0.52
23:BB:1720:U:O2'	23:BB:1721:G:H5'	2.10	0.52
23:BB:1838:C:H4'	23:BB:1839:G:C8	2.45	0.52
23:BB:231:A:H3'	23:BB:232:G:H8	1.75	0.52
23:BB:2852:G:H2'	23:BB:2853:C:O4'	2.10	0.52
25:BC:132:ARG:NH2	25:BC:169:ALA:HA	2.25	0.52
25:BC:20:ASN:O	25:BC:23:LEU:HB2	2.10	0.52
48:BG:97:VAL:HB	48:BG:124:CYS:HB2	1.91	0.52
48:BG:84:LYS:O	48:BG:85:LYS:HG2	2.10	0.52
40:BH:5:LEU:HD13	40:BH:13:GLY:N	2.23	0.52
24:BI:129:GLU:CB	24:BI:133:ARG:HH12	2.17	0.52
42:BN:58:ASP:O	42:BN:62:ASN:HB2	2.09	0.52
49:BR:49:ILE:HG12	49:BR:53:PHE:CA	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BS:24:ILE:CD1	45:BS:36:LEU:HD21	2.40	0.52
35:BV:80:HIS:HB3	35:BV:83:LYS:O	2.09	0.52
1:CA:10:A:OP2	4:CE:130:THR:HB	2.09	0.52
1:CA:1476:A:O2'	1:CA:1477:U:H5'	2.09	0.52
1:CA:50:A:N6	1:CA:361:G:H4'	2.25	0.52
1:CA:565:U:H3'	1:CA:566:G:H2'	1.92	0.52
1:CA:675:A:H2'	1:CA:676:A:H8	1.74	0.52
1:CA:824:G:O2'	1:CA:825:A:H5'	2.09	0.52
1:CA:757:U:O2'	1:CA:879:C:H1'	2.09	0.52
18:CB:16:GLY:HA2	18:CB:40:ILE:HG13	1.92	0.52
18:CB:217:ALA:O	18:CB:220:VAL:HB	2.10	0.52
2:CC:149:LYS:HG2	2:CC:200:TRP:CE3	2.43	0.52
2:CC:51:VAL:C	2:CC:114:LEU:HD21	2.29	0.52
4:CE:55:VAL:N	4:CE:56:PRO:HD2	2.25	0.52
7:CH:17:GLN:OE1	7:CH:69:ALA:HB1	2.10	0.52
9:CJ:59:LYS:HG3	9:CJ:60:ASP:N	2.24	0.52
20:CO:29:VAL:HG11	20:CO:67:LEU:HD21	1.90	0.52
13:CP:48:GLU:CD	13:CP:49:GLY:H	2.13	0.52
17:CT:73:ARG:HB3	17:CT:73:ARG:HH11	1.74	0.52
23:DB:1533:C:O2'	23:DB:1534:U:H5'	2.10	0.52
23:DB:2617:U:C2'	23:DB:2618:G:H5'	2.39	0.52
23:DB:871:U:H2'	23:DB:872:U:C6	2.45	0.52
23:DB:877:A:H2'	23:DB:877:A:N3	2.24	0.52
23:DB:876:C:H3'	23:DB:877:A:H4'	1.91	0.52
23:DB:942:G:O2'	23:DB:943:A:H5'	2.09	0.52
25:DC:202:ARG:NH1	25:DC:204:LEU:HD11	2.25	0.52
25:DC:43:ASN:ND2	25:DC:44:ASN:H	2.07	0.52
23:DB:2621:G:P	26:DD:124:ARG:HH22	2.33	0.52
47:DF:91:ARG:HD3	47:DF:91:ARG:N	2.25	0.52
47:DF:91:ARG:C	47:DF:95:MET:HB2	2.30	0.52
27:DK:64:ARG:O	27:DK:82:ASN:HA	2.10	0.52
42:DN:9:GLN:O	42:DN:11:ASN:N	2.43	0.52
42:DN:55:ALA:HA	42:DN:80:PHE:CD1	2.45	0.52
43:DO:116:GLN:N	43:DO:116:GLN:HE21	2.08	0.52
43:DO:28:VAL:O	43:DO:28:VAL:HG13	2.10	0.52
44:DQ:105:PHE:O	44:DQ:108:LEU:HB2	2.10	0.52
50:DT:11:LEU:HD21	50:DT:46:ALA:HB1	1.92	0.52
1:AA:1376:U:H2'	1:AA:1377:A:H8	1.75	0.52
1:AA:13:U:O2	1:AA:914:A:H3'	2.10	0.52
1:AA:241:G:O2'	1:AA:242:G:H5'	2.10	0.52
1:AA:399:G:H2'	1:AA:400:C:C6	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:729:A:H2'	1:AA:730:G:C8	2.45	0.52
1:AA:784:A:H2'	1:AA:785:G:C8	2.45	0.52
1:AA:796:C:O2'	1:AA:797:C:H5'	2.10	0.52
1:AA:892:A:H2'	1:AA:893:C:C6	2.45	0.52
18:AB:148:GLY:HA2	18:AB:151:LYS:HD2	1.92	0.52
18:AB:205:ALA:O	18:AB:209:VAL:HG22	2.10	0.52
2:AC:129:PHE:HB2	2:AC:133:MET:HE3	1.92	0.52
3:AD:144:ILE:HD13	3:AD:154:VAL:HG11	1.91	0.52
4:AE:37:VAL:HG12	4:AE:47:PHE:HB2	1.91	0.52
4:AE:56:PRO:HG2	4:AE:57:ALA:H	1.75	0.52
4:AE:63:MET:O	4:AE:66:ALA:HB3	2.10	0.52
8:AI:25:GLY:HA3	8:AI:57:VAL:O	2.10	0.52
20:AO:35:GLN:O	20:AO:39:LEU:N	2.42	0.52
20:AO:78:TYR:OH	20:AO:88:ARG:HD2	2.09	0.52
36:B2:45:SER:O	36:B2:46:LYS:CB	2.55	0.52
23:BB:1151:A:H2'	23:BB:1152:C:H6	1.74	0.52
23:BB:1535:A:H2	23:BB:1538:G:H1'	1.75	0.52
23:BB:1947:C:H2'	23:BB:1948:G:C8	2.40	0.52
23:BB:2281:A:H4'	23:BB:2389:G:N2	2.25	0.52
23:BB:2369:A:H2'	23:BB:2370:G:H8	1.74	0.52
23:BB:2586:U:H2'	23:BB:2587:A:H8	1.74	0.52
23:BB:2636:C:O2'	23:BB:2637:U:H5'	2.10	0.52
23:BB:2784:U:O2'	23:BB:2785:C:H5'	2.10	0.52
23:BB:2786:U:H4'	26:BD:67:HIS:HA	1.91	0.52
23:BB:693:A:O2'	23:BB:694:U:H5'	2.09	0.52
23:BB:962:G:H2'	23:BB:963:U:H6	1.74	0.52
25:BC:2:VAL:HG23	25:BC:3:VAL:N	2.20	0.52
26:BD:154:LYS:H	26:BD:154:LYS:HD3	1.75	0.52
23:BB:2636:C:H4'	26:BD:81:GLU:OE2	2.10	0.52
47:BF:116:LEU:HG	47:BF:117:SER:H	1.73	0.52
48:BG:153:PRO:CG	48:BG:162:ARG:HB3	2.36	0.52
24:BI:17:ALA:O	24:BI:18:ASN:CB	2.58	0.52
27:BK:25:LEU:HB2	27:BK:38:ILE:HG12	1.90	0.52
27:BK:64:ARG:HD2	27:BK:102:PRO:O	2.10	0.52
42:BN:51:LEU:O	42:BN:54:LEU:HB3	2.10	0.52
45:BS:55:ILE:HD12	45:BS:107:VAL:HG21	1.92	0.52
50:BT:57:VAL:HG13	50:BT:58:VAL:N	2.24	0.52
50:BT:69:ARG:CA	50:BT:69:ARG:HE	2.21	0.52
46:BU:35:VAL:CG2	46:BU:38:ILE:HG21	2.40	0.52
1:CA:1306:A:O2'	1:CA:1307:U:H5'	2.10	0.52
1:CA:22:G:H2'	1:CA:23:C:H6	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:386:C:C2'	1:CA:387:U:H5'	2.40	0.52
18:CB:82:ALA:CB	18:CB:213:LEU:HD13	2.40	0.52
3:CD:55:ARG:NE	3:CD:55:ARG:HA	2.24	0.52
7:CH:29:SER:OG	7:CH:32:LYS:HG3	2.10	0.52
9:CJ:32:THR:O	9:CJ:80:THR:HG21	2.10	0.52
23:DB:1544:A:H2'	23:DB:1545:A:C8	2.45	0.52
23:DB:1930:G:H2'	23:DB:1968:G:N1	2.25	0.52
23:DB:2015:A:C2	31:D0:2:VAL:HG22	2.45	0.52
23:DB:2037:A:H2'	23:DB:2038:G:C8	2.45	0.52
23:DB:208:C:H2'	23:DB:209:C:C6	2.45	0.52
23:DB:2210:U:N3	23:DB:2212:A:N7	2.57	0.52
23:DB:234:U:H2'	23:DB:235:U:H6	1.74	0.52
23:DB:2357:G:N2	23:DB:2359:C:H3'	2.25	0.52
23:DB:2466:C:O2'	23:DB:2467:C:H5'	2.09	0.52
23:DB:287:G:H2'	23:DB:288:U:O4'	2.10	0.52
23:DB:409:G:H2'	23:DB:410:G:C8	2.44	0.52
23:DB:4:U:H2'	23:DB:5:A:C8	2.45	0.52
23:DB:909:A:H2'	23:DB:912:C:H5	1.73	0.52
25:DC:71:ASP:HB3	25:DC:118:GLY:HA2	1.92	0.52
26:DD:136:ASN:HD21	26:DD:139:SER:C	2.12	0.52
26:DD:79:LEU:HD22	26:DD:79:LEU:N	2.24	0.52
47:DF:136:ILE:HG22	47:DF:142:TYR:HB3	1.92	0.52
47:DF:28:PRO:O	47:DF:168:LEU:HD11	2.10	0.52
47:DF:40:GLY:O	47:DF:41:GLU:HB2	2.10	0.52
47:DF:45:ASP:O	47:DF:46:LYS:HE2	2.10	0.52
23:DB:2529:G:H5'	48:DG:174:LYS:HG3	1.92	0.52
48:DG:26:LYS:HG2	48:DG:27:GLY:N	2.25	0.52
40:DH:5:LEU:HD22	40:DH:12:LEU:O	2.10	0.52
38:DM:63:ILE:HD12	38:DM:63:ILE:H	1.75	0.52
44:DQ:17:LEU:HD13	44:DQ:30:VAL:O	2.09	0.52
52:DW:81:ILE:O	52:DW:81:ILE:HG13	2.10	0.52
51:DZ:33:LEU:H	51:DZ:52:SER:CB	2.23	0.52
1:AA:1343:G:O2'	1:AA:1344:C:H5'	2.09	0.52
1:AA:1527:U:O2'	1:AA:1528:U:H5'	2.09	0.52
1:AA:384:G:H2'	1:AA:385:C:C6	2.44	0.52
1:AA:580:C:H2'	1:AA:581:G:C8	2.45	0.52
1:AA:895:G:H2'	1:AA:896:C:C6	2.44	0.52
1:AA:975:A:N6	1:AA:1366:C:O2'	2.42	0.52
4:AE:136:VAL:HG13	4:AE:137:ARG:H	1.74	0.52
4:AE:55:VAL:N	4:AE:56:PRO:HD2	2.25	0.52
11:AL:31:GLY:HA3	11:AL:54:VAL:HG13	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:552:U:H4'	11:AL:82:ARG:HG2	1.92	0.52
12:AM:9:PRO:C	12:AM:44:ILE:HD13	2.30	0.52
21:AN:12:ARG:HG3	21:AN:53:ASP:OD2	2.09	0.52
23:BB:1103:A:H3'	23:BB:1104:C:H6	1.75	0.52
23:BB:1120:G:O2'	23:BB:1121:C:H5'	2.10	0.52
23:BB:1520:U:H2'	23:BB:1521:G:O4'	2.09	0.52
23:BB:1723:G:N7	23:BB:1737:G:N2	2.54	0.52
23:BB:1742:U:O2'	23:BB:1743:G:H5'	2.10	0.52
23:BB:1782:U:H2'	23:BB:1783:A:H5'	1.91	0.52
23:BB:1878:G:H2'	23:BB:1879:C:C6	2.45	0.52
23:BB:2251:G:H2'	23:BB:2252:G:C8	2.45	0.52
23:BB:233:A:H61	23:BB:428:A:H61	1.58	0.52
23:BB:2344:U:H4'	23:BB:2345:G:OP1	2.08	0.52
23:BB:2530:A:H3'	48:BG:156:TYR:OH	2.09	0.52
23:BB:2611:C:O2'	23:BB:2612:C:H5'	2.09	0.52
23:BB:871:U:H1'	23:BB:907:G:N1	2.25	0.52
23:BB:920:A:H2'	23:BB:921:C:H6	1.74	0.52
26:BD:117:GLY:HA2	26:BD:164:GLN:OE1	2.10	0.52
29:BE:148:ILE:HG13	29:BE:167:VAL:CG2	2.40	0.52
29:BE:165:HIS:C	29:BE:167:VAL:H	2.13	0.52
29:BE:192:ALA:O	29:BE:196:VAL:HG23	2.10	0.52
29:BE:29:HIS:NE2	37:BL:8:PRO:HG3	2.25	0.52
47:BF:105:ILE:C	47:BF:108:PRO:HD2	2.30	0.52
40:BH:73:ASN:HB2	40:BH:141:LYS:HB2	1.92	0.52
40:BH:79:THR:HG22	40:BH:145:ASN:OD1	2.10	0.52
40:BH:14:SER:HB2	40:BH:17:ASP:CB	2.39	0.52
40:BH:4:ILE:O	40:BH:36:ALA:HA	2.10	0.52
40:BH:47:PHE:O	40:BH:50:ARG:HB2	2.10	0.52
24:BI:23:VAL:HG23	24:BI:24:GLY:N	2.25	0.52
41:BJ:106:LYS:HE3	41:BJ:106:LYS:O	2.10	0.52
38:BM:35:ALA:HB3	38:BM:99:GLY:N	2.24	0.52
52:BW:24:ARG:HD3	52:BW:65:LYS:HG2	1.90	0.52
1:CA:216:U:H2'	1:CA:217:C:C6	2.45	0.52
1:CA:253:A:H2'	1:CA:254:G:H8	1.73	0.52
1:CA:746:A:H2'	1:CA:747:A:H8	1.75	0.52
1:CA:889:A:H5'	1:CA:891:U:H1'	1.90	0.52
18:CB:101:THR:HG22	18:CB:174:GLU:OE1	2.09	0.52
8:CI:16:ALA:HB2	8:CI:66:VAL:CG2	2.35	0.52
8:CI:46:VAL:HA	8:CI:49:GLN:HG3	1.91	0.52
21:CN:3:GLN:OE1	21:CN:6:LYS:HG3	2.10	0.52
16:CS:16:LYS:HD2	16:CS:17:LYS:NZ	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:CT:53:MET:HA	17:CT:56:ILE:CD1	2.39	0.52
36:D2:3:ARG:HE	36:D2:4:THR:HG22	1.74	0.52
23:DB:1405:U:H2'	23:DB:1406:U:C6	2.45	0.52
23:DB:1506:U:H2'	23:DB:1507:C:C6	2.45	0.52
23:DB:1550:C:H2'	23:DB:1551:A:H8	1.75	0.52
23:DB:1825:U:H2'	23:DB:1826:G:C8	2.45	0.52
23:DB:1842:G:H2'	23:DB:1843:C:H6	1.73	0.52
23:DB:235:U:H2'	23:DB:236:C:C6	2.44	0.52
23:DB:62:U:H3'	23:DB:63:A:C8	2.45	0.52
23:DB:796:C:H2'	23:DB:797:G:H8	1.75	0.52
25:DC:260:LYS:C	25:DC:262:THR:H	2.13	0.52
26:DD:40:LEU:HA	26:DD:45:TYR:H	1.74	0.52
29:DE:109:LEU:HA	29:DE:112:LEU:HB2	1.91	0.52
29:DE:192:ALA:O	29:DE:196:VAL:HG23	2.10	0.52
29:DE:48:THR:C	29:DE:50:ALA:H	2.11	0.52
23:DB:2305:U:H5''	47:DF:130:GLY:HA3	1.92	0.52
40:DH:116:ARG:N	40:DH:131:SER:O	2.38	0.52
24:DI:54:ILE:C	24:DI:54:ILE:HD13	2.29	0.52
41:DJ:34:ARG:NH1	41:DJ:39:LYS:HG2	2.25	0.52
38:DM:102:LEU:HD22	38:DM:102:LEU:H	1.74	0.52
28:DP:31:VAL:HG11	28:DP:38:ARG:HG2	1.91	0.52
26:DD:10:GLY:HA2	28:DP:4:ILE:HD11	1.92	0.52
46:DU:26:ASN:H	46:DU:26:ASN:HD22	1.57	0.52
22:DA:83:G:OP1	30:DY:16:LEU:HD23	2.10	0.52
1:AA:1078:U:H4'	4:AE:137:ARG:NH1	2.24	0.52
1:AA:1144:G:N2	1:AA:1146:A:H62	2.08	0.52
1:AA:1243:C:H2'	1:AA:1244:G:C8	2.45	0.52
1:AA:1434:A:H2'	1:AA:1435:G:C8	2.44	0.52
1:AA:731:G:OP1	1:AA:766:A:H1'	2.10	0.52
6:AG:52:ARG:NH2	6:AG:120:ALA:HB1	2.23	0.52
6:AG:74:VAL:HG21	6:AG:85:GLN:NE2	2.24	0.52
11:AL:51:VAL:HG12	11:AL:52:CYS:N	2.19	0.52
11:AL:79:ILE:HD12	11:AL:96:THR:HG22	1.91	0.52
16:AS:63:ASP:O	16:AS:66:VAL:HG13	2.10	0.52
17:AT:43:LYS:N	17:AT:43:LYS:HZ2	2.08	0.52
23:BB:1141:U:H5''	41:BJ:27:ARG:NH2	2.25	0.52
23:BB:2264:C:H41	52:BW:11:ASN:ND2	2.08	0.52
23:BB:2345:G:H5'	23:BB:2347:C:O4'	2.10	0.52
23:BB:267:C:H2'	23:BB:268:C:C6	2.44	0.52
23:BB:467:G:O2'	23:BB:468:G:H5'	2.10	0.52
23:BB:62:U:H3'	23:BB:63:A:C8	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:862:G:H2'	23:BB:863:A:C8	2.45	0.52
25:BC:270:ARG:HB3	25:BC:270:ARG:NH1	2.25	0.52
29:BE:58:LYS:CD	29:BE:58:LYS:N	2.72	0.52
23:BB:1061:U:H5'	24:BI:9:LYS:NZ	2.25	0.52
27:BK:24:VAL:HA	27:BK:39:ILE:CD1	2.40	0.52
42:BN:62:ASN:N	42:BN:62:ASN:HD22	2.06	0.52
27:BK:76:VAL:H	28:BP:72:VAL:HG23	1.74	0.52
50:BT:29:THR:HA	50:BT:86:THR:CA	2.33	0.52
1:CA:1307:U:H2'	1:CA:1308:U:C6	2.45	0.52
1:CA:1480:A:H2'	1:CA:1481:U:H6	1.75	0.52
1:CA:624:C:O2'	1:CA:625:U:H5'	2.10	0.52
1:CA:692:U:O2	1:CA:694:A:H5''	2.10	0.52
3:CD:197:HIS:ND1	3:CD:198:LEU:N	2.58	0.52
5:CF:3:HIS:CG	5:CF:92:THR:HG23	2.45	0.52
6:CG:55:LYS:HB2	6:CG:59:GLU:OE2	2.09	0.52
16:CS:35:ARG:HB2	16:CS:71:GLY:CA	2.40	0.52
17:CT:47:GLN:HG2	17:CT:82:ILE:HD13	1.91	0.52
23:DB:1178:C:H2'	23:DB:1179:G:C8	2.45	0.52
23:DB:1930:G:H2'	23:DB:1968:G:C6	2.45	0.52
23:DB:570:G:H2'	23:DB:2030:A:C8	2.45	0.52
23:DB:2590:A:H2'	23:DB:2591:C:H6	1.74	0.52
23:DB:666:A:H4'	37:DL:48:ARG:HD2	1.92	0.52
23:DB:721:A:H2'	23:DB:722:A:H8	1.75	0.52
25:DC:158:GLY:N	25:DC:194:VAL:HG13	2.25	0.52
29:DE:157:LEU:HG	29:DE:169:VAL:HG11	1.92	0.52
23:DB:659:G:N2	29:DE:30:GLN:HE22	2.04	0.52
47:DF:3:LEU:O	47:DF:3:LEU:HD13	2.09	0.52
50:DT:40:LYS:HA	50:DT:43:ILE:CG2	2.40	0.52
52:DW:46:ALA:HB2	52:DW:78:PHE:CD1	2.43	0.52
1:AA:1463:U:H2'	1:AA:1464:U:H6	1.76	0.51
1:AA:642:A:H2'	1:AA:643:C:C6	2.44	0.51
1:AA:645:G:H2'	1:AA:646:G:H8	1.75	0.51
1:AA:953:G:H2'	1:AA:954:G:O4'	2.10	0.51
18:AB:81:ASP:O	18:AB:84:LEU:HD23	2.11	0.51
2:AC:9:ILE:HG23	2:AC:10:ARG:HG3	1.92	0.51
3:AD:123:MET:SD	3:AD:127:ARG:N	2.83	0.51
3:AD:153:ARG:HG3	3:AD:154:VAL:N	2.25	0.51
1:AA:546:A:P	3:AD:68:GLU:HB3	2.50	0.51
5:AF:3:HIS:ND1	5:AF:95:ALA:N	2.57	0.51
6:AG:50:ALA:HB1	6:AG:55:LYS:O	2.10	0.51
8:AI:38:PHE:O	8:AI:44:ARG:HB3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:54:VAL:O	11:AL:61:GLU:HA	2.10	0.51
12:AM:95:PRO:N	12:AM:108:ARG:HG2	2.26	0.51
19:AU:42:THR:HB	19:AU:46:ARG:HE	1.75	0.51
23:BB:1010:A:H1'	23:BB:1153:C:O4'	2.10	0.51
23:BB:108:G:H2'	23:BB:109:C:O4'	2.10	0.51
23:BB:1151:A:H2'	23:BB:1152:C:C6	2.45	0.51
23:BB:1279:G:H4'	42:BN:31:HIS:CD2	2.44	0.51
23:BB:1439:A:N7	23:BB:1440:U:N1	2.59	0.51
23:BB:1599:U:H2'	23:BB:1600:C:H6	1.74	0.51
23:BB:721:A:H2'	23:BB:722:A:H8	1.75	0.51
25:BC:91:ALA:HB2	25:BC:105:ALA:HB2	1.90	0.51
25:BC:71:ASP:HB3	25:BC:118:GLY:HA2	1.92	0.51
48:BG:43:LYS:HA	48:BG:43:LYS:HE2	1.92	0.51
48:BG:85:LYS:HB2	48:BG:164:ALA:HB2	1.91	0.51
40:BH:121:VAL:HB	40:BH:128:HIS:CE1	2.44	0.51
40:BH:79:THR:HA	40:BH:144:VAL:HA	1.91	0.51
37:BL:125:LEU:HB2	37:BL:143:GLU:OE2	2.09	0.51
28:BP:60:VAL:O	28:BP:70:GLU:HA	2.10	0.51
44:BQ:79:ILE:HG23	44:BQ:80:ASN:N	2.25	0.51
44:BQ:89:ILE:HG21	49:BR:11:GLN:NE2	2.25	0.51
45:BS:74:ILE:HD12	45:BS:104:THR:O	2.10	0.51
45:BS:36:LEU:HD11	45:BS:47:VAL:HB	1.91	0.51
50:BT:11:LEU:HD21	50:BT:46:ALA:HB1	1.91	0.51
35:BV:14:LYS:CE	35:BV:18:ARG:HH21	2.22	0.51
35:BV:5:ASN:O	35:BV:6:ALA:HB2	2.10	0.51
30:BY:58:GLU:H	30:BY:58:GLU:CD	2.13	0.51
1:CA:1209:C:O2'	1:CA:1210:C:H5'	2.10	0.51
1:CA:123:U:H5''	1:CA:311:C:O2'	2.10	0.51
1:CA:143:A:H2	1:CA:220:G:H22	1.57	0.51
1:CA:399:G:H2'	1:CA:400:C:C6	2.45	0.51
2:CC:22:PHE:HZ	9:CJ:11:LYS:HZ2	1.57	0.51
4:CE:44:ARG:HA	4:CE:71:ILE:O	2.09	0.51
8:CI:44:ARG:HB2	8:CI:48:ARG:NH2	2.26	0.51
10:CK:80:ASN:CA	10:CK:105:ARG:HB3	2.38	0.51
11:CL:33:CYS:HA	11:CL:54:VAL:HA	1.92	0.51
1:CA:1216:A:H5''	21:CN:4:SER:CB	2.39	0.51
14:CQ:68:LYS:HG2	14:CQ:69:THR:HG23	1.91	0.51
22:DA:26:C:H3'	22:DA:27:C:C6	2.44	0.51
23:DB:1083:U:H2'	23:DB:1085:A:OP2	2.10	0.51
23:DB:1439:A:C6	23:DB:1552:A:N7	2.78	0.51
23:DB:2322:A:H3'	23:DB:2323:G:C8	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2369:A:H2'	23:DB:2370:G:H8	1.74	0.51
23:DB:471:A:H2'	23:DB:472:A:O4'	2.10	0.51
23:DB:571:U:N3	23:DB:575:A:N7	2.58	0.51
23:DB:680:C:H2'	23:DB:681:G:H8	1.76	0.51
23:DB:78:U:H2'	23:DB:79:C:C6	2.45	0.51
25:DC:92:LEU:HD11	25:DC:100:ARG:HB2	1.92	0.51
23:DB:1826:G:P	25:DC:221:GLY:H	2.34	0.51
25:DC:244:VAL:HG12	25:DC:250:GLN:N	2.26	0.51
25:DC:52:HIS:CD2	25:DC:52:HIS:N	2.79	0.51
47:DF:135:ILE:HG13	47:DF:137:PHE:CD2	2.44	0.51
22:DA:43:C:H4'	47:DF:91:ARG:HG3	1.92	0.51
48:DG:84:LYS:O	48:DG:85:LYS:HG2	2.10	0.51
40:DH:135:HIS:HB3	40:DH:138:VAL:CG2	2.41	0.51
24:DI:79:LEU:HD11	24:DI:131:THR:OG1	2.10	0.51
37:DL:132:ARG:HA	37:DL:135:ILE:HG22	1.90	0.51
37:DL:9:ALA:HB3	37:DL:12:SER:OG	2.10	0.51
23:DB:910:A:H62	38:DM:12:MET:HA	1.74	0.51
28:DP:63:ILE:O	28:DP:63:ILE:HG22	2.10	0.51
41:DJ:44:TYR:CD2	44:DQ:59:LEU:HD21	2.44	0.51
50:DT:30:ILE:HG23	50:DT:85:VAL:CG2	2.40	0.51
1:AA:1391:U:H2'	1:AA:1392:G:N7	2.26	0.51
1:AA:1432:G:H1'	1:AA:1468:A:H61	1.75	0.51
1:AA:618:C:N3	1:AA:622:A:N6	2.58	0.51
1:AA:663:A:O2'	1:AA:664:G:H5'	2.10	0.51
1:AA:81:A:H2'	1:AA:81:A:N3	2.24	0.51
2:AC:155:ARG:HH12	2:AC:160:GLU:HA	1.75	0.51
3:AD:144:ILE:HG22	3:AD:145:ARG:N	2.24	0.51
3:AD:54:LEU:CD1	3:AD:55:ARG:HH21	2.24	0.51
3:AD:88:ASN:O	3:AD:92:LEU:HD23	2.09	0.51
13:AP:48:GLU:CD	13:AP:49:GLY:H	2.14	0.51
23:BB:1175:A:H3'	23:BB:1176:U:O4'	2.10	0.51
23:BB:1475:G:H1'	23:BB:1476:U:H5	1.75	0.51
23:BB:1533:C:O2'	23:BB:1534:U:H5'	2.11	0.51
23:BB:1677:A:H2'	23:BB:1678:A:C8	2.45	0.51
23:BB:1872:A:O5'	23:BB:1872:A:H8	1.92	0.51
23:BB:1999:C:O2'	23:BB:2000:C:H5'	2.10	0.51
23:BB:1133:A:N6	23:BB:2025:C:O2'	2.43	0.51
23:BB:2047:C:H2'	23:BB:2048:G:C8	2.45	0.51
23:BB:2665:A:O2'	23:BB:2666:C:H5'	2.09	0.51
23:BB:393:C:O2'	23:BB:394:C:H5'	2.11	0.51
23:BB:575:A:O2'	23:BB:576:U:H5'	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:736:C:H2'	23:BB:737:C:H6	1.75	0.51
25:BC:260:LYS:C	25:BC:262:THR:H	2.13	0.51
47:BF:136:ILE:HG22	47:BF:142:TYR:HB3	1.93	0.51
47:BF:107:VAL:HG21	47:BF:175:PRO:HG3	1.93	0.51
48:BG:67:ALA:O	48:BG:71:LEU:HD23	2.10	0.51
24:BI:52:LEU:HD21	24:BI:81:LYS:HZ2	1.75	0.51
41:BJ:114:LEU:O	41:BJ:118:MET:HG3	2.10	0.51
41:BJ:4:PHE:CG	41:BJ:5:THR:N	2.78	0.51
38:BM:90:GLU:HA	38:BM:90:GLU:OE1	2.11	0.51
43:BO:67:ASN:HB3	43:BO:70:ALA:CB	2.40	0.51
50:BT:7:LEU:O	50:BT:7:LEU:HD13	2.11	0.51
52:BW:13:ARG:HG3	52:BW:14:ASP:H	1.75	0.51
1:CA:1090:U:H2'	1:CA:1091:U:C6	2.45	0.51
1:CA:1434:A:H2'	1:CA:1435:G:C8	2.45	0.51
1:CA:977:A:H2'	1:CA:978:A:C5'	2.39	0.51
2:CC:123:LEU:HA	2:CC:127:VAL:HG22	1.91	0.51
2:CC:143:LEU:O	2:CC:143:LEU:HD13	2.10	0.51
2:CC:38:VAL:HG23	2:CC:39:ARG:N	2.24	0.51
4:CE:85:LYS:HG3	4:CE:93:VAL:O	2.11	0.51
6:CG:48:THR:O	6:CG:52:ARG:HG3	2.11	0.51
9:CJ:56:HIS:O	9:CJ:57:VAL:HG12	2.10	0.51
9:CJ:93:ALA:HB3	9:CJ:96:VAL:HG22	1.90	0.51
10:CK:20:ALA:O	10:CK:22:ILE:HG13	2.10	0.51
11:CL:110:LYS:O	11:CL:113:ARG:HG3	2.10	0.51
11:CL:49:ARG:CB	11:CL:89:LEU:HD11	2.39	0.51
13:CP:71:VAL:HG13	13:CP:72:ALA:N	2.25	0.51
14:CQ:32:ILE:HG23	14:CQ:33:TYR:CD2	2.45	0.51
17:CT:75:LYS:O	17:CT:78:LEU:HB2	2.10	0.51
23:DB:1275:A:O2'	23:DB:1276:A:H5'	2.09	0.51
23:DB:1723:G:H3'	23:DB:1724:G:C8	2.39	0.51
23:DB:2199:A:H5'	23:DB:2200:C:OP2	2.10	0.51
23:DB:2803:G:H2'	23:DB:2804:U:H6	1.76	0.51
23:DB:688:U:H2'	23:DB:689:A:H8	1.75	0.51
23:DB:693:A:H2'	23:DB:694:U:C6	2.45	0.51
23:DB:836:G:H2'	23:DB:837:C:C6	2.45	0.51
25:DC:20:ASN:HD21	25:DC:22:GLU:HG3	1.74	0.51
47:DF:115:GLY:HA3	47:DF:177:ARG:HD2	1.90	0.51
24:DI:32:VAL:HG22	24:DI:60:VAL:CG2	2.40	0.51
23:DB:1203:U:O4'	37:DL:3:LEU:HD12	2.10	0.51
44:DQ:63:ARG:HH22	44:DQ:96:ASP:CA	2.23	0.51
1:AA:1003:G:N2	1:AA:1005:A:H5'	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1120:C:H2'	1:AA:1121:U:H6	1.76	0.51
1:AA:1241:G:H2'	1:AA:1242:G:H8	1.74	0.51
1:AA:944:G:N1	1:AA:1338:G:OP2	2.44	0.51
1:AA:1366:C:O2'	1:AA:1367:C:H5'	2.10	0.51
1:AA:251:G:H22	1:AA:272:C:N4	2.09	0.51
1:AA:651:C:H2'	1:AA:652:U:C6	2.44	0.51
18:AB:86:CYS:C	18:AB:88:GLN:H	2.13	0.51
2:AC:69:THR:HG21	2:AC:75:VAL:HG21	1.92	0.51
1:AA:10:A:OP2	4:AE:130:THR:HB	2.10	0.51
5:AF:3:HIS:N	5:AF:3:HIS:CD2	2.78	0.51
6:AG:87:PRO:HG3	6:AG:144:ALA:HA	1.93	0.51
20:AO:56:LEU:O	20:AO:59:MET:HG2	2.09	0.51
16:AS:52:ASN:CG	16:AS:53:GLY:N	2.63	0.51
17:AT:49:ALA:O	17:AT:52:GLU:HB3	2.11	0.51
22:BA:43:C:C2'	22:BA:44:G:H5''	2.39	0.51
23:BB:1199:U:H2'	23:BB:1200:C:C6	2.45	0.51
23:BB:1355:G:O2'	23:BB:1356:G:H5'	2.09	0.51
23:BB:1405:U:H2'	23:BB:1406:U:C6	2.45	0.51
23:BB:140:C:C5'	23:BB:141:G:H21	2.23	0.51
23:BB:1494:A:H2'	23:BB:1495:A:C8	2.45	0.51
23:BB:1689:A:H2'	23:BB:1690:A:C8	2.46	0.51
23:BB:2359:C:H2'	23:BB:2360:G:C8	2.44	0.51
23:BB:2869:G:H2'	23:BB:2870:C:C6	2.45	0.51
23:BB:967:U:H2'	23:BB:968:C:C6	2.46	0.51
47:BF:3:LEU:HD13	47:BF:3:LEU:O	2.10	0.51
47:BF:91:ARG:HD3	47:BF:91:ARG:N	2.25	0.51
48:BG:166:GLU:CD	48:BG:166:GLU:N	2.63	0.51
48:BG:24:THR:C	48:BG:25:ILE:HD12	2.30	0.51
40:BH:5:LEU:HD22	40:BH:12:LEU:O	2.10	0.51
40:BH:96:THR:HG22	40:BH:115:VAL:HG11	1.91	0.51
24:BI:100:ILE:O	24:BI:139:VAL:HA	2.10	0.51
1:CA:332:G:O2'	1:CA:333:U:H5'	2.10	0.51
1:CA:747:A:C4	1:CA:748:G:H1'	2.46	0.51
18:CB:80:LYS:HE2	18:CB:81:ASP:CG	2.31	0.51
3:CD:144:ILE:HG22	3:CD:145:ARG:N	2.26	0.51
3:CD:169:TRP:CD2	3:CD:185:PRO:HB3	2.46	0.51
4:CE:146:MET:HG3	4:CE:146:MET:O	2.10	0.51
6:CG:26:VAL:CB	6:CG:42:VAL:HG21	2.41	0.51
20:CO:74:ASP:OD1	20:CO:76:ALA:HB3	2.11	0.51
7:CH:81:GLY:O	14:CQ:35:LYS:HD3	2.09	0.51
1:CA:673:A:H1'	15:CR:63:TYR:HE1	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:CT:43:LYS:HB3	17:CT:86:ALA:CB	2.39	0.51
23:DB:1309:G:H4'	36:D2:7:PRO:HB2	1.93	0.51
23:DB:1808:A:O2'	23:DB:1809:A:H5'	2.11	0.51
23:DB:1826:G:H2'	23:DB:1827:U:H6	1.75	0.51
23:DB:2835:A:N6	23:DB:2878:U:H2'	2.25	0.51
23:DB:57:C:H2'	23:DB:58:G:H8	1.76	0.51
23:DB:64:A:H2'	23:DB:65:U:H6	1.75	0.51
48:DG:85:LYS:HB2	48:DG:164:ALA:HB2	1.92	0.51
24:DI:126:ARG:HH11	24:DI:126:ARG:HB3	1.75	0.51
24:DI:63:ASP:O	24:DI:64:ARG:HB2	2.09	0.51
37:DL:109:LYS:HB2	37:DL:111:ILE:HD13	1.92	0.51
42:DN:101:GLY:CA	42:DN:110:MET:H	2.15	0.51
46:DU:86:PHE:HE1	46:DU:88:ASP:HB2	1.75	0.51
1:AA:103:U:H2'	1:AA:104:G:O4'	2.10	0.51
1:AA:1102:A:H2'	1:AA:1103:C:C6	2.45	0.51
1:AA:476:U:H2'	1:AA:477:C:C6	2.45	0.51
1:AA:555:U:H2'	1:AA:556:C:H6	1.76	0.51
1:AA:580:C:H2'	1:AA:581:G:O4'	2.10	0.51
1:AA:824:G:O2'	1:AA:825:A:H5'	2.11	0.51
2:AC:149:LYS:HG3	2:AC:168:ARG:HG3	1.92	0.51
3:AD:13:ARG:HG3	3:AD:55:ARG:NH1	2.26	0.51
4:AE:136:VAL:HG13	4:AE:137:ARG:N	2.25	0.51
4:AE:97:PRO:HA	4:AE:122:VAL:HG12	1.91	0.51
6:AG:78:ARG:CZ	6:AG:78:ARG:HB3	2.40	0.51
1:AA:1129:C:H5''	8:AI:17:ARG:HH22	1.76	0.51
10:AK:51:PHE:HZ	10:AK:61:ALA:HA	1.75	0.51
10:AK:85:VAL:O	10:AK:111:ASP:HA	2.10	0.51
12:AM:89:ARG:HH22	12:AM:101:THR:HG21	1.74	0.51
17:AT:73:ARG:HB3	17:AT:73:ARG:HH11	1.76	0.51
34:B3:31:ILE:HD11	34:B3:34:LYS:CD	2.39	0.51
23:BB:1429:G:H2'	23:BB:1430:G:C8	2.45	0.51
23:BB:144:A:H2'	23:BB:145:C:C6	2.46	0.51
23:BB:1513:U:O2'	23:BB:1514:G:H5'	2.10	0.51
23:BB:1680:U:O2	23:BB:1763:G:H3'	2.09	0.51
23:BB:1773:A:H2'	23:BB:1774:C:O4'	2.10	0.51
23:BB:1846:G:N2	23:BB:1848:A:N6	2.58	0.51
23:BB:1909:C:H2'	23:BB:1910:G:H8	1.75	0.51
23:BB:2250:G:N2	38:BM:82:MET:HB3	2.25	0.51
23:BB:2617:U:O2'	23:BB:2618:G:H5'	2.10	0.51
23:BB:693:A:H2'	23:BB:694:U:C6	2.45	0.51
23:BB:903:C:H2'	23:BB:904:G:H8	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:99:U:O2	23:BB:99:U:O4'	2.27	0.51
25:BC:70:LYS:HD3	25:BC:95:TYR:CE1	2.46	0.51
26:BD:123:LYS:HD2	26:BD:165:MET:HE1	1.92	0.51
29:BE:148:ILE:HG22	29:BE:148:ILE:O	2.10	0.51
23:BB:674:G:H5''	29:BE:71:GLY:CA	2.40	0.51
47:BF:100:GLU:O	47:BF:104:THR:HB	2.11	0.51
24:BI:125:THR:O	24:BI:129:GLU:HG3	2.10	0.51
27:BK:19:VAL:HB	27:BK:41:ILE:HD11	1.90	0.51
46:BU:85:ARG:HD3	46:BU:86:PHE:N	2.26	0.51
52:BW:39:GLN:HG2	52:BW:40:ARG:H	1.75	0.51
52:BW:69:GLU:HG3	52:BW:80:SER:OG	2.10	0.51
1:CA:1299:A:C8	1:CA:1301:U:H1'	2.44	0.51
1:CA:384:G:H2'	1:CA:385:C:C6	2.45	0.51
1:CA:731:G:OP1	1:CA:766:A:H1'	2.11	0.51
1:CA:789:U:H2'	1:CA:791:G:OP2	2.11	0.51
1:CA:818:G:C3'	1:CA:819:A:H5''	2.40	0.51
5:CF:53:LYS:N	5:CF:53:LYS:NZ	2.59	0.51
11:CL:7:VAL:HG22	14:CQ:33:TYR:HD1	1.76	0.51
21:CN:24:ALA:O	21:CN:27:LYS:HG2	2.10	0.51
22:DA:16:G:O2'	22:DA:17:C:H5'	2.11	0.51
23:DB:1654:A:H2'	23:DB:1655:A:H8	1.76	0.51
23:DB:1714:U:H3'	23:DB:1715:G:C5'	2.40	0.51
23:DB:176:A:O2'	23:DB:177:G:H5'	2.10	0.51
23:DB:1773:A:H2'	23:DB:1774:C:O4'	2.11	0.51
23:DB:1830:C:H2'	23:DB:1831:G:C8	2.46	0.51
23:DB:1908:C:O2'	23:DB:1909:C:H5'	2.11	0.51
23:DB:1925:C:H2'	23:DB:1926:U:H5''	1.93	0.51
23:DB:1947:C:H2'	23:DB:1948:G:C8	2.42	0.51
23:DB:2773:C:H2'	23:DB:2774:C:H6	1.75	0.51
23:DB:360:U:H2'	23:DB:361:G:C8	2.45	0.51
23:DB:521:U:H2'	23:DB:522:A:C8	2.46	0.51
29:DE:14:VAL:HG12	29:DE:197:GLU:OE1	2.09	0.51
24:DI:78:LEU:HD13	24:DI:108:ILE:HG23	1.93	0.51
38:DM:57:VAL:O	38:DM:57:VAL:HG12	2.11	0.51
42:DN:87:PHE:HB3	42:DN:90:ARG:CB	2.39	0.51
49:DR:14:VAL:HG23	49:DR:18:GLN:HG3	1.92	0.51
46:DU:78:LYS:CD	46:DU:79:ALA:H	2.23	0.51
35:DV:51:GLN:NE2	35:DV:79:ARG:HH22	2.08	0.51
35:DV:5:ASN:O	35:DV:6:ALA:HB2	2.10	0.51
39:DX:23:ARG:HA	39:DX:27:ASN:H	1.76	0.51
30:DY:10:ARG:HB3	30:DY:10:ARG:NH2	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1151:A:O2'	1:AA:1152:A:H8	1.94	0.51
1:AA:1240:U:H3'	1:AA:1241:G:H5'	1.91	0.51
1:AA:1340:A:O2'	1:AA:1341:U:H5'	2.11	0.51
1:AA:451:A:H1'	1:AA:452:A:C8	2.46	0.51
1:AA:861:G:H2'	1:AA:862:C:C6	2.44	0.51
18:AB:186:VAL:HG21	18:AB:192:PRO:HB3	1.93	0.51
2:AC:51:VAL:HA	2:AC:69:THR:HA	1.92	0.51
3:AD:8:LEU:O	3:AD:12:ARG:HG2	2.10	0.51
13:AP:71:VAL:HG13	13:AP:72:ALA:N	2.26	0.51
19:AU:16:ARG:HH12	19:AU:19:LYS:NZ	2.08	0.51
23:BB:1430:G:H2'	23:BB:1431:A:H8	1.76	0.51
23:BB:1825:U:H2'	23:BB:1826:G:H8	1.75	0.51
23:BB:1830:C:H2'	23:BB:1831:G:C8	2.46	0.51
23:BB:2261:C:O2'	23:BB:2262:U:H5'	2.09	0.51
23:BB:2322:A:H3'	23:BB:2323:G:C8	2.46	0.51
23:BB:978:G:O2'	23:BB:979:A:H5'	2.11	0.51
25:BC:166:ARG:HA	25:BC:171:VAL:HG22	1.92	0.51
26:BD:79:LEU:N	26:BD:79:LEU:HD22	2.25	0.51
29:BE:126:VAL:HG13	29:BE:127:GLU:N	2.26	0.51
47:BF:2:LYS:HB2	47:BF:100:GLU:OE2	2.10	0.51
47:BF:155:ILE:HG22	47:BF:157:THR:H	1.75	0.51
22:BA:43:C:C1'	47:BF:91:ARG:HD2	2.40	0.51
40:BH:72:ILE:HG23	40:BH:140:ALA:HB1	1.91	0.51
40:BH:96:THR:HA	40:BH:115:VAL:HG11	1.92	0.51
41:BJ:34:ARG:NH1	41:BJ:39:LYS:HG2	2.26	0.51
41:BJ:58:ASN:HA	41:BJ:127:GLY:CA	2.40	0.51
31:B0:53:VAL:HB	42:BN:118:ARG:HH22	1.75	0.51
42:BN:47:VAL:O	42:BN:50:PRO:HD2	2.10	0.51
42:BN:87:PHE:HB3	42:BN:90:ARG:CB	2.41	0.51
43:BO:30:ARG:HG2	43:BO:31:THR:N	2.26	0.51
35:BV:41:GLU:C	35:BV:42:LEU:HD23	2.30	0.51
23:BB:2352:A:N1	52:BW:30:VAL:HG11	2.24	0.51
52:BW:55:ASP:C	52:BW:57:THR:H	2.13	0.51
1:CA:8:A:C6	3:CD:205:LYS:HB2	2.45	0.51
1:CA:811:C:O2'	1:CA:901:A:N1	2.42	0.51
2:CC:172:VAL:O	2:CC:174:LEU:HD12	2.11	0.51
4:CE:96:GLN:HB3	4:CE:123:LEU:CD1	2.41	0.51
5:CF:69:GLU:O	5:CF:73:GLU:HG2	2.10	0.51
12:CM:64:VAL:HA	12:CM:68:LEU:CD1	2.41	0.51
1:CA:980:C:H4'	21:CN:12:ARG:NH2	2.25	0.51
13:CP:5:ARG:HH22	13:CP:24:SER:HA	1.70	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1026:G:H2'	23:DB:1027:A:C8	2.46	0.51
23:DB:1060:U:O2	23:DB:1088:A:C8	2.64	0.51
23:DB:1771:C:H2'	23:DB:1772:A:H8	1.76	0.51
23:DB:2023:C:O2'	23:DB:2024:G:H5'	2.10	0.51
23:DB:2417:C:O2'	23:DB:2418:A:H5'	2.11	0.51
23:DB:2656:U:H2'	23:DB:2657:A:H8	1.75	0.51
25:DC:12:ARG:HA	25:DC:15:VAL:HG23	1.93	0.51
26:DD:124:ARG:HD3	26:DD:163:GLY:O	2.10	0.51
47:DF:57:ALA:HB2	47:DF:64:PRO:HG2	1.92	0.51
48:DG:33:THR:HA	48:DG:34:ARG:NH1	2.25	0.51
24:DI:85:ILE:CD1	24:DI:137:LEU:HD21	2.40	0.51
38:DM:40:ARG:HB2	38:DM:93:VAL:HG22	1.91	0.51
42:DN:90:ARG:HB3	42:DN:94:TYR:HE1	1.76	0.51
52:DW:23:LYS:O	52:DW:66:VAL:HB	2.11	0.51
1:AA:1010:U:O2'	1:AA:1011:C:H5'	2.10	0.51
1:AA:1070:U:H2'	1:AA:1071:C:H6	1.73	0.51
1:AA:1142:G:H3'	1:AA:1143:G:H8	1.76	0.51
1:AA:218:U:H2'	1:AA:219:U:C6	2.46	0.51
1:AA:434:U:H3'	1:AA:435:A:H8	1.75	0.51
1:AA:499:A:H1'	1:AA:500:G:C8	2.45	0.51
1:AA:57:G:H2'	1:AA:58:C:H6	1.76	0.51
1:AA:62:U:H2'	1:AA:63:C:C6	2.46	0.51
1:AA:714:G:H2'	1:AA:715:A:C8	2.46	0.51
2:AC:39:ARG:CZ	2:AC:56:ILE:HD11	2.40	0.51
4:AE:113:VAL:O	4:AE:116:VAL:HG22	2.11	0.51
11:AL:81:ILE:HA	11:AL:95:HIS:O	2.11	0.51
14:AQ:26:ARG:HG2	14:AQ:39:ARG:O	2.10	0.51
23:BB:1275:A:O2'	23:BB:1276:A:H5'	2.11	0.51
23:BB:138:U:H3'	23:BB:140:C:O2	2.11	0.51
23:BB:151:C:H2'	23:BB:152:A:C8	2.46	0.51
23:BB:1714:U:H3'	23:BB:1715:G:C5'	2.40	0.51
23:BB:1917:U:H2'	23:BB:1918:A:C8	2.46	0.51
23:BB:1951:U:O2	23:BB:1953:A:H3'	2.11	0.51
23:BB:2337:G:N3	23:BB:2337:G:H2'	2.26	0.51
23:BB:2803:G:H2'	23:BB:2804:U:H6	1.75	0.51
23:BB:500:G:N2	23:BB:502:A:H3'	2.25	0.51
23:BB:959:A:O2'	23:BB:960:A:H5'	2.11	0.51
48:BG:89:VAL:HG12	48:BG:90:GLY:N	2.26	0.51
40:BH:68:ARG:HG3	40:BH:134:VAL:HB	1.93	0.51
40:BH:97:ARG:HA	40:BH:112:LYS:CB	2.41	0.51
27:BK:71:ARG:CB	27:BK:72:PRO:HD2	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BM:40:ARG:HB2	38:BM:93:VAL:HG22	1.92	0.51
44:BQ:17:LEU:HD13	44:BQ:30:VAL:O	2.11	0.51
1:CA:1125:U:O2'	1:CA:1126:U:H2'	2.10	0.51
1:CA:1298:U:C3'	1:CA:1299:A:H5'	2.40	0.51
1:CA:1349:A:OP1	8:CI:121:ARG:HB2	2.09	0.51
1:CA:1451:U:H5''	1:CA:1452:C:H5	1.75	0.51
1:CA:1465:A:H2'	1:CA:1466:C:C6	2.45	0.51
1:CA:148:G:H2'	1:CA:149:A:O4'	2.10	0.51
1:CA:184:G:N1	1:CA:194:C:N4	2.59	0.51
1:CA:279:A:H5'	1:CA:281:G:C5'	2.40	0.51
7:CH:25:THR:O	7:CH:26:MET:HB3	2.11	0.51
9:CJ:59:LYS:HB2	9:CJ:62:ARG:HH22	1.74	0.51
10:CK:83:VAL:CG2	10:CK:109:ILE:HG12	2.41	0.51
10:CK:126:ARG:HE	10:CK:126:ARG:HA	1.76	0.51
11:CL:106:VAL:HA	11:CL:107:LYS:NZ	2.25	0.51
31:D0:38:LEU:HB3	31:D0:41:HIS:CD2	2.45	0.51
22:DA:43:C:C2'	22:DA:44:G:H5''	2.40	0.51
22:DA:76:G:H2'	22:DA:77:U:H6	1.74	0.51
23:DB:1151:A:H2'	23:DB:1152:C:C6	2.45	0.51
23:DB:1447:C:H2'	23:DB:1448:G:C8	2.46	0.51
1:CA:1409:C:O2	23:DB:1913:A:N6	2.44	0.51
23:DB:528:A:N1	23:DB:2042:A:H2'	2.25	0.51
23:DB:2144:G:N3	23:DB:2146:C:H5'	2.26	0.51
23:DB:2443:C:H2'	23:DB:2444:G:H8	1.76	0.51
23:DB:2579:C:O5'	23:DB:2579:C:H6	1.94	0.51
23:DB:267:C:H2'	23:DB:268:C:C6	2.45	0.51
23:DB:282:A:O2'	23:DB:283:G:H5'	2.10	0.51
23:DB:2852:G:H2'	23:DB:2853:C:O4'	2.10	0.51
23:DB:340:A:H2'	23:DB:341:C:H5'	1.92	0.51
23:DB:693:A:O2'	23:DB:694:U:H5'	2.10	0.51
29:DE:77:ILE:HG13	29:DE:78:TRP:HE3	1.76	0.51
47:DF:3:LEU:HD11	47:DF:172:PHE:CE1	2.45	0.51
48:DG:89:VAL:HG12	48:DG:90:GLY:N	2.26	0.51
27:DK:16:ALA:HB3	27:DK:47:ILE:HG13	1.92	0.51
37:DL:100:ILE:O	37:DL:100:ILE:HG12	2.10	0.51
28:DP:91:VAL:O	28:DP:92:ARG:HB3	2.10	0.51
44:DQ:79:ILE:HG23	44:DQ:80:ASN:N	2.25	0.51
44:DQ:91:ARG:HD3	49:DR:11:GLN:CD	2.31	0.51
45:DS:48:LYS:O	45:DS:51:LEU:HB3	2.10	0.51
50:DT:35:ALA:O	50:DT:81:LYS:HB3	2.10	0.51
46:DU:43:LYS:HD3	46:DU:44:HIS:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DW:19:ARG:HD3	52:DW:36:ILE:HD11	1.93	0.51
39:DX:46:VAL:O	39:DX:50:VAL:HG23	2.11	0.51
1:AA:1015:G:H1'	1:AA:1218:C:O2'	2.11	0.51
1:AA:1120:C:H2'	1:AA:1121:U:C6	2.45	0.51
1:AA:1360:A:H2'	1:AA:1361:G:H5'	1.91	0.51
1:AA:648:A:H2'	1:AA:649:A:C8	2.46	0.51
1:AA:658:C:O2'	1:AA:659:U:H5'	2.10	0.51
1:AA:735:C:H2'	1:AA:736:C:H6	1.76	0.51
1:AA:735:C:O2'	1:AA:736:C:H5'	2.10	0.51
18:AB:125:PHE:C	18:AB:127:LYS:H	2.14	0.51
2:AC:126:ARG:HH22	2:AC:190:THR:HG22	1.75	0.51
2:AC:38:VAL:HG23	2:AC:39:ARG:H	1.75	0.51
8:AI:54:VAL:O	8:AI:59:LYS:HD2	2.10	0.51
11:AL:37:TYR:O	11:AL:38:THR:HG23	2.11	0.51
12:AM:64:VAL:HB	12:AM:65:GLU:OE2	2.11	0.51
23:BB:130:C:O2'	23:BB:131:A:H5'	2.11	0.51
23:BB:2065:C:H2'	23:BB:2066:C:H6	1.76	0.51
23:BB:235:U:H2'	23:BB:236:C:C6	2.46	0.51
23:BB:252:G:O2'	23:BB:253:C:H5'	2.11	0.51
23:BB:2590:A:H2'	23:BB:2591:C:C6	2.45	0.51
23:BB:37:C:H4'	23:BB:451:U:OP1	2.10	0.51
23:BB:599:A:O2'	23:BB:600:G:H5'	2.10	0.51
23:BB:64:A:H2'	23:BB:65:U:C6	2.45	0.51
23:BB:939:G:O2'	23:BB:940:G:H5'	2.10	0.51
26:BD:5:VAL:HG23	26:BD:32:ASN:ND2	2.25	0.51
29:BE:14:VAL:HG12	29:BE:197:GLU:OE1	2.10	0.51
47:BF:107:VAL:N	47:BF:108:PRO:CD	2.74	0.51
47:BF:7:TYR:O	47:BF:12:VAL:HG23	2.11	0.51
47:BF:78:ILE:HA	47:BF:82:TYR:CD1	2.45	0.51
40:BH:57:LYS:HG3	40:BH:58:LEU:N	2.25	0.51
37:BL:132:ARG:HA	37:BL:135:ILE:CG2	2.41	0.51
37:BL:19:LEU:CD2	37:BL:27:LEU:HD23	2.40	0.51
44:BQ:105:PHE:O	44:BQ:108:LEU:HB2	2.11	0.51
49:BR:66:HIS:ND1	49:BR:94:THR:HG22	2.26	0.51
45:BS:37:THR:HG22	45:BS:48:LYS:HE3	1.92	0.51
46:BU:49:PRO:HA	46:BU:53:GLN:HE21	1.74	0.51
1:CA:1051:C:H2'	1:CA:1052:U:C6	2.46	0.51
1:CA:1118:U:H2'	1:CA:1119:C:C6	2.45	0.51
1:CA:1468:A:O2'	1:CA:1469:C:H5'	2.10	0.51
1:CA:33:A:H2'	1:CA:34:C:H6	1.74	0.51
1:CA:493:A:H3'	1:CA:494:G:C8	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:502:A:H4'	1:CA:550:G:H4'	1.92	0.51
1:CA:555:U:H2'	1:CA:556:C:C6	2.45	0.51
1:CA:868:C:H2'	1:CA:869:G:O4'	2.11	0.51
2:CC:64:ARG:HG3	2:CC:101:ASN:HD21	1.75	0.51
3:CD:156:ALA:O	3:CD:159:GLU:HB3	2.11	0.51
8:CI:20:ILE:CG1	8:CI:62:LEU:HD12	2.41	0.51
9:CJ:42:LEU:HD22	9:CJ:73:LEU:HD23	1.93	0.51
11:CL:54:VAL:O	11:CL:61:GLU:HA	2.10	0.51
12:CM:15:VAL:O	12:CM:19:THR:HG23	2.10	0.51
21:CN:5:MET:HB3	21:CN:62:ARG:HH12	1.76	0.51
17:CT:49:ALA:O	17:CT:52:GLU:HB3	2.10	0.51
33:D1:36:LYS:CB	33:D1:47:ILE:HA	2.40	0.51
23:DB:1169:A:O2'	23:DB:1170:C:H5'	2.11	0.51
23:DB:1672:A:H2'	23:DB:1673:G:C8	2.45	0.51
23:DB:1946:U:H2'	23:DB:1947:C:C6	2.45	0.51
23:DB:2052:A:H4'	26:DD:148:GLN:O	2.11	0.51
23:DB:2305:U:H2'	23:DB:2306:C:C6	2.46	0.51
23:DB:2800:A:H2'	23:DB:2801:G:C4'	2.41	0.51
23:DB:2834:G:H1'	23:DB:2883:A:N6	2.26	0.51
23:DB:458:G:H22	23:DB:469:G:H2'	1.76	0.51
23:DB:607:U:O4	23:DB:620:G:H5''	2.11	0.51
29:DE:46:GLN:HB3	29:DE:86:ALA:CA	2.41	0.51
27:DK:107:LEU:N	27:DK:107:LEU:HD12	2.21	0.51
37:DL:14:LYS:O	37:DL:16:GLY:N	2.44	0.51
49:DR:1:MET:HA	49:DR:42:ALA:HB3	1.92	0.51
1:AA:1007:U:H2'	1:AA:1008:U:C6	2.46	0.51
1:AA:1020:G:C2'	1:AA:1021:A:H5'	2.40	0.51
1:AA:1285:A:O2'	1:AA:1286:U:OP2	2.25	0.51
1:AA:1310:G:H2'	1:AA:1311:A:O4'	2.11	0.51
1:AA:282:A:H2'	1:AA:282:A:N3	2.26	0.51
1:AA:33:A:H2'	1:AA:34:C:H6	1.74	0.51
1:AA:403:C:O2'	1:AA:404:G:H5'	2.09	0.51
1:AA:531:U:H6	1:AA:531:U:H5'	1.75	0.51
1:AA:600:A:H2'	1:AA:601:G:H8	1.76	0.51
1:AA:624:C:O2'	1:AA:625:U:H5'	2.10	0.51
1:AA:93:U:H3'	1:AA:94:G:H5''	1.93	0.51
1:AA:987:G:H2'	1:AA:988:G:C8	2.44	0.51
18:AB:16:GLY:HA2	18:AB:40:ILE:HG13	1.92	0.51
3:AD:145:ARG:HB3	3:AD:147:LYS:HG3	1.92	0.51
4:AE:45:VAL:HG11	4:AE:117:ALA:HB2	1.92	0.51
10:AK:70:ALA:HA	10:AK:73:VAL:CG2	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:86:VAL:CG2	11:AL:89:LEU:HB2	2.40	0.51
12:AM:92:ARG:HE	12:AM:92:ARG:HA	1.75	0.51
13:AP:6:LEU:HB3	13:AP:17:TYR:HB3	1.93	0.51
1:AA:1319:A:H3'	16:AS:2:ARG:CA	2.41	0.51
23:BB:1178:C:H2'	23:BB:1179:G:C8	2.46	0.51
23:BB:1281:G:H2'	23:BB:1282:U:C6	2.46	0.51
23:BB:1287:A:H3'	23:BB:1288:G:H21	1.75	0.51
23:BB:142:A:O2'	50:BT:2:ILE:HG21	2.11	0.51
23:BB:1567:G:H1'	23:BB:1568:G:C6	2.46	0.51
23:BB:1841:U:H2'	23:BB:1842:G:H8	1.76	0.51
25:BC:91:ALA:CB	25:BC:105:ALA:HB2	2.41	0.51
23:BB:2653:U:O2'	48:BG:109:SER:HB2	2.10	0.51
40:BH:67:ALA:N	40:BH:138:VAL:HG21	2.26	0.51
38:BM:21:ALA:HB3	38:BM:99:GLY:O	2.11	0.51
44:BQ:90:ASP:OD1	49:BR:39:LEU:HD12	2.11	0.51
50:BT:39:THR:O	50:BT:40:LYS:HB3	2.11	0.51
46:BU:11:ILE:O	46:BU:12:VAL:HB	2.10	0.51
46:BU:41:VAL:O	46:BU:42:LYS:HB2	2.11	0.51
1:CA:1058:G:OP1	2:CC:198:LYS:HE2	2.11	0.51
1:CA:137:U:H2'	1:CA:138:G:H8	1.74	0.51
1:CA:299:G:H2'	1:CA:300:A:C8	2.46	0.51
1:CA:692:U:C2	1:CA:694:A:H5''	2.46	0.51
1:CA:699:C:C2'	1:CA:700:G:H5''	2.40	0.51
1:CA:736:C:H2'	1:CA:737:C:H6	1.75	0.51
1:CA:82:G:H2'	1:CA:83:C:O4'	2.11	0.51
18:CB:114:LYS:HZ3	18:CB:151:LYS:HD2	1.75	0.51
3:CD:52:VAL:HG12	3:CD:198:LEU:HD11	1.93	0.51
3:CD:88:ASN:O	3:CD:92:LEU:HD23	2.11	0.51
5:CF:22:ILE:O	5:CF:26:THR:HG23	2.11	0.51
11:CL:37:TYR:O	11:CL:38:THR:HG23	2.11	0.51
21:CN:51:PRO:HG2	21:CN:52:ARG:H	1.76	0.51
1:CA:981:U:C4'	21:CN:60:ARG:HD2	2.35	0.51
23:DB:2419:U:H5''	33:D1:21:THR:HG21	1.93	0.51
23:DB:108:G:O2'	23:DB:109:C:H5'	2.10	0.51
23:DB:135:U:H2'	23:DB:136:G:H8	1.75	0.51
23:DB:1655:A:H2'	23:DB:1656:C:O4'	2.10	0.51
23:DB:2457:U:O2'	23:DB:2458:G:H5'	2.11	0.51
23:DB:2880:C:O2'	23:DB:2881:U:H5'	2.11	0.51
23:DB:441:U:H2'	23:DB:442:G:H8	1.75	0.51
23:DB:500:G:N2	23:DB:502:A:H3'	2.25	0.51
23:DB:543:G:H2'	23:DB:545:U:H5'	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DE:158:PHE:HA	29:DE:169:VAL:CG2	2.40	0.51
29:DE:28:VAL:O	29:DE:32:VAL:HG22	2.11	0.51
47:DF:62:GLN:NE2	47:DF:91:ARG:HE	2.08	0.51
48:DG:62:ALA:O	48:DG:66:THR:HG22	2.09	0.51
41:DJ:20:ALA:HB2	41:DJ:28:LEU:HD13	1.93	0.51
27:DK:21:CYS:HB2	27:DK:39:ILE:HG21	1.93	0.51
38:DM:32:GLY:HA3	38:DM:103:TYR:O	2.11	0.51
28:DP:88:ARG:HB2	28:DP:112:ARG:CZ	2.41	0.51
28:DP:19:PHE:CE1	28:DP:46:VAL:HG11	2.46	0.51
45:DS:33:LEU:O	45:DS:37:THR:HG23	2.11	0.51
35:DV:38:LEU:HG	35:DV:40:ILE:HG23	1.93	0.51
35:DV:80:HIS:HB3	35:DV:83:LYS:O	2.10	0.51
52:DW:13:ARG:HG3	52:DW:14:ASP:H	1.75	0.51
1:AA:1090:U:H2'	1:AA:1091:U:C6	2.46	0.51
1:AA:1314:C:H2'	1:AA:1315:U:C6	2.46	0.51
1:AA:1508:A:H2'	1:AA:1509:C:H6	1.75	0.51
1:AA:187:G:N2	1:AA:189:A:H3'	2.25	0.51
1:AA:313:A:H2'	1:AA:314:C:H6	1.76	0.51
1:AA:476:U:H2'	1:AA:477:C:H6	1.76	0.51
1:AA:599:C:O2'	1:AA:600:A:H5'	2.10	0.51
1:AA:951:G:O2'	1:AA:952:U:H5'	2.11	0.51
18:AB:110:ILE:C	18:AB:112:ARG:H	2.14	0.51
10:AK:22:ILE:HG21	10:AK:95:THR:HG21	1.92	0.51
31:B0:38:LEU:HB3	31:B0:41:HIS:CD2	2.46	0.51
34:B3:22:LYS:HG3	34:B3:47:ALA:O	2.11	0.51
23:BB:100:U:O2'	23:BB:101:A:O5'	2.26	0.51
23:BB:1458:U:O2'	23:BB:1459:G:H4'	2.09	0.51
23:BB:1595:C:O2'	23:BB:1596:A:H5'	2.10	0.51
23:BB:1676:A:H2'	23:BB:1677:A:O4'	2.11	0.51
23:BB:1825:U:H2'	23:BB:1826:G:C8	2.46	0.51
23:BB:1827:U:O2'	23:BB:1828:G:H5'	2.11	0.51
23:BB:2047:C:H2'	23:BB:2048:G:H8	1.75	0.51
23:BB:19:A:H2'	23:BB:20:C:C6	2.46	0.51
23:BB:64:A:H2'	23:BB:65:U:H6	1.76	0.51
23:BB:721:A:H2'	23:BB:722:A:C8	2.46	0.51
25:BC:189:ALA:O	25:BC:190:THR:C	2.49	0.51
26:BD:61:THR:OG1	26:BD:64:GLU:HB2	2.11	0.51
48:BG:26:LYS:HG2	48:BG:27:GLY:N	2.24	0.51
24:BI:77:VAL:HA	24:BI:80:LYS:CE	2.38	0.51
27:BK:41:ILE:HG13	27:BK:42:THR:H	1.75	0.51
37:BL:65:GLY:O	37:BL:66:PHE:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BM:54:THR:C	38:BM:56:ALA:H	2.14	0.51
44:BQ:93:ILE:HG23	44:BQ:94:LEU:H	1.76	0.51
49:BR:38:VAL:HG13	49:BR:54:VAL:HG12	1.93	0.51
1:CA:1207:G:O2'	1:CA:1208:C:H5'	2.11	0.51
1:CA:735:C:H2'	1:CA:736:C:H6	1.75	0.51
1:CA:821:G:H2'	1:CA:822:U:C6	2.46	0.51
2:CC:172:VAL:HG11	2:CC:200:TRP:HB3	1.91	0.51
3:CD:129:VAL:HG12	3:CD:130:ASN:N	2.26	0.51
3:CD:40:HIS:O	3:CD:43:ARG:HG2	2.11	0.51
11:CL:49:ARG:HH12	11:CL:88:ASP:HB3	1.76	0.51
11:CL:52:CYS:SG	11:CL:66:ILE:HD11	2.50	0.51
12:CM:74:MET:SD	12:CM:77:LYS:HD2	2.51	0.51
12:CM:78:ARG:HB3	12:CM:78:ARG:CZ	2.40	0.51
12:CM:89:ARG:HD3	12:CM:94:LEU:HB2	1.92	0.51
23:DB:1205:A:N7	29:DE:165:HIS:CG	2.78	0.51
23:DB:138:U:H5''	23:DB:139:U:OP1	2.11	0.51
23:DB:1774:C:O2	23:DB:1774:C:C2'	2.59	0.51
23:DB:1973:G:O2'	23:DB:1974:C:H5'	2.11	0.51
23:DB:2037:A:H2'	23:DB:2038:G:H8	1.76	0.51
26:DD:106:LYS:HD3	26:DD:106:LYS:N	2.26	0.51
26:DD:32:ASN:HB3	26:DD:50:VAL:HG21	1.91	0.51
47:DF:32:LYS:H	47:DF:95:MET:HE1	1.75	0.51
48:DG:24:THR:C	48:DG:25:ILE:HD12	2.32	0.51
40:DH:114:GLU:HG3	40:DH:133:GLN:HB2	1.93	0.51
40:DH:84:ALA:HB3	40:DH:148:ALA:HB1	1.92	0.51
40:DH:69:ALA:O	40:DH:72:ILE:HG13	2.11	0.51
40:DH:84:ALA:C	40:DH:91:PHE:HE1	2.14	0.51
41:DJ:4:PHE:CG	41:DJ:5:THR:N	2.78	0.51
41:DJ:65:THR:HG23	41:DJ:66:GLY:N	2.26	0.51
42:DN:58:ASP:O	42:DN:62:ASN:HB2	2.10	0.51
42:DN:92:GLY:O	42:DN:94:TYR:N	2.44	0.51
50:DT:43:ILE:CG2	50:DT:58:VAL:HG21	2.41	0.51
50:DT:29:THR:CG2	50:DT:86:THR:HG22	2.41	0.51
46:DU:28:LEU:HB2	46:DU:32:LYS:O	2.10	0.51
1:AA:1073:U:O2'	1:AA:1074:G:H5'	2.09	0.51
1:AA:1378:C:N3	1:AA:1379:G:H1'	2.26	0.51
1:AA:803:G:H2'	1:AA:804:U:H6	1.73	0.51
1:AA:929:G:O2'	1:AA:930:C:H5'	2.11	0.51
1:AA:976:G:N2	1:AA:1362:A:H3'	2.25	0.51
18:AB:11:ALA:C	18:AB:13:VAL:H	2.13	0.51
18:AB:221:ARG:CB	18:AB:221:ARG:HH11	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:148:ILE:O	2:AC:168:ARG:HG2	2.10	0.51
3:AD:169:TRP:CD2	3:AD:185:PRO:HB3	2.45	0.51
3:AD:199:ILE:HG13	3:AD:200:VAL:N	2.25	0.51
5:AF:6:ILE:HD12	5:AF:7:VAL:N	2.25	0.51
8:AI:21:LYS:HG2	8:AI:22:PRO:HD2	1.92	0.51
1:AA:1152:A:H4'	9:AJ:15:HIS:CD2	2.45	0.51
10:AK:126:ARG:HA	10:AK:126:ARG:HE	1.76	0.51
20:AO:33:THR:HG23	20:AO:63:ARG:NH1	2.24	0.51
16:AS:38:THR:HG23	16:AS:69:LYS:HE2	1.92	0.51
23:BB:1106:G:H2'	23:BB:1107:G:C8	2.44	0.51
23:BB:1117:C:H2'	23:BB:1118:C:H6	1.75	0.51
23:BB:1558:C:H4'	23:BB:1559:U:C5'	2.36	0.51
23:BB:1710:G:O2'	23:BB:1711:A:H5'	2.10	0.51
23:BB:570:G:H2'	23:BB:2030:A:C8	2.46	0.51
23:BB:2394:C:H2'	23:BB:2395:C:C6	2.46	0.51
23:BB:39:G:O2'	23:BB:40:U:H5'	2.10	0.51
23:BB:593:U:H2'	23:BB:594:U:H6	1.76	0.51
23:BB:666:A:H2'	23:BB:667:U:C6	2.46	0.51
23:BB:5:A:H2'	23:BB:6:A:H8	1.72	0.51
23:BB:866:A:H61	23:BB:913:U:C4'	2.24	0.51
26:BD:51:THR:HG23	26:BD:78:GLY:O	2.11	0.51
29:BE:46:GLN:HB3	29:BE:86:ALA:CA	2.40	0.51
37:BL:142:ILE:HD12	37:BL:142:ILE:N	2.26	0.51
37:BL:51:GLU:OE1	37:BL:57:LEU:HB2	2.11	0.51
44:BQ:94:LEU:C	44:BQ:96:ASP:N	2.64	0.51
52:BW:17:ALA:O	52:BW:18:LYS:HD2	2.10	0.51
39:BX:20:ASN:O	39:BX:24:GLU:HB3	2.11	0.51
1:CA:1004:A:H2'	1:CA:1005:A:O4'	2.11	0.51
1:CA:1271:A:H5'	1:CA:1314:C:C5'	2.40	0.51
1:CA:475:C:O2'	1:CA:476:U:H5'	2.11	0.51
5:CF:84:VAL:HG22	5:CF:85:ILE:H	1.74	0.51
6:CG:12:LEU:HD13	6:CG:13:PRO:CD	2.41	0.51
6:CG:12:LEU:HD13	6:CG:13:PRO:HD2	1.92	0.51
1:CA:1351:U:O4'	6:CG:32:ASP:HB3	2.11	0.51
8:CI:11:ARG:O	8:CI:11:ARG:HG3	2.11	0.51
8:CI:29:ILE:HG23	8:CI:64:ILE:HB	1.93	0.51
20:CO:36:ILE:HD11	20:CO:59:MET:HG3	1.93	0.51
20:CO:81:LEU:C	20:CO:81:LEU:HD23	2.32	0.51
34:D3:31:ILE:CD1	34:D3:34:LYS:HD3	2.41	0.51
23:DB:1000:A:H2'	23:DB:1001:A:H8	1.76	0.51
23:DB:1019:U:H2'	23:DB:1020:A:H8	1.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:172:A:H2'	23:DB:173:A:H8	1.75	0.51
23:DB:222:A:N6	23:DB:232:G:H1'	2.26	0.51
23:DB:2531:A:N3	23:DB:2531:A:H2'	2.26	0.51
23:DB:2694:G:H2'	23:DB:2695:U:C6	2.46	0.51
23:DB:68:G:O2'	23:DB:69:C:H5'	2.11	0.51
25:DC:140:VAL:HG12	25:DC:141:HIS:N	2.22	0.51
25:DC:204:LEU:HD23	25:DC:209:ALA:HB3	1.91	0.51
25:DC:66:PHE:HB2	25:DC:150:GLY:O	2.11	0.51
47:DF:134:GLN:HB3	47:DF:149:ARG:CB	2.41	0.51
27:DK:24:VAL:HA	27:DK:39:ILE:CD1	2.41	0.51
37:DL:136:GLU:C	37:DL:138:ALA:H	2.15	0.51
38:DM:33:LEU:HD22	38:DM:128:THR:CB	2.41	0.51
42:DN:9:GLN:HA	42:DN:17:ARG:HD3	1.92	0.51
28:DP:24:THR:OG1	28:DP:86:LYS:HB3	2.10	0.51
44:DQ:93:ILE:HG23	44:DQ:94:LEU:H	1.76	0.51
45:DS:24:ILE:HD11	45:DS:36:LEU:HD21	1.91	0.51
1:AA:977:A:N1	1:AA:1224:U:OP1	2.45	0.50
1:AA:1299:A:H2'	1:AA:1301:U:C6	2.46	0.50
1:AA:336:A:O2'	1:AA:337:G:H5'	2.10	0.50
1:AA:91:U:H2'	1:AA:92:U:C5	2.46	0.50
7:AH:37:ASN:O	7:AH:40:LYS:HB3	2.11	0.50
31:B0:27:LEU:HD22	31:B0:38:LEU:HA	1.92	0.50
22:BA:26:C:H3'	22:BA:27:C:C6	2.45	0.50
23:BB:1431:A:H2'	23:BB:1432:G:C8	2.45	0.50
23:BB:1826:G:H2'	23:BB:1827:U:H6	1.76	0.50
23:BB:2109:U:O2	23:BB:2109:U:H2'	2.10	0.50
23:BB:2352:A:C6	52:BW:30:VAL:HG11	2.47	0.50
23:BB:2355:G:H4'	52:BW:20:LEU:HD13	1.93	0.50
23:BB:2773:C:H2'	23:BB:2774:C:H6	1.76	0.50
23:BB:350:G:O2'	23:BB:351:C:H5'	2.11	0.50
23:BB:569:U:H2'	23:BB:570:G:O4'	2.10	0.50
23:BB:1826:G:P	25:BC:221:GLY:H	2.34	0.50
25:BC:180:MET:HB2	25:BC:268:ARG:HB3	1.93	0.50
26:BD:117:GLY:HA2	26:BD:164:GLN:CD	2.32	0.50
24:BI:29:GLN:HA	24:BI:29:GLN:HE21	1.76	0.50
27:BK:116:ILE:HG13	27:BK:117:SER:N	2.26	0.50
52:BW:50:VAL:HG23	52:BW:61:LYS:CE	2.41	0.50
1:CA:1078:U:H2'	1:CA:1079:G:O4'	2.10	0.50
1:CA:370:C:H2'	1:CA:371:A:H8	1.75	0.50
1:CA:470:C:H2'	1:CA:471:U:H6	1.77	0.50
1:CA:94:G:H4'	1:CA:95:C:O5'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:105:VAL:HA	2:CC:106:ARG:CZ	2.42	0.50
2:CC:153:SER:HG	2:CC:196:GLY:H	1.57	0.50
2:CC:113:LYS:HA	2:CC:184:ASN:ND2	2.26	0.50
1:CA:795:C:H5''	10:CK:127:ARG:HH21	1.76	0.50
17:CT:38:ILE:HD11	17:CT:82:ILE:HA	1.93	0.50
23:DB:1050:A:O2'	23:DB:2752:C:H1'	2.11	0.50
23:DB:1431:A:H2'	23:DB:1432:G:H8	1.76	0.50
23:DB:1459:G:H3'	23:DB:1460:U:C5'	2.40	0.50
23:DB:1545:A:H2'	23:DB:1546:G:O4'	2.12	0.50
23:DB:171:U:H2'	23:DB:172:A:H8	1.74	0.50
23:DB:1720:U:O2'	23:DB:1721:G:H5'	2.11	0.50
23:DB:2852:G:H2'	23:DB:2853:C:H6	1.76	0.50
23:DB:4:U:H2'	23:DB:5:A:H8	1.76	0.50
23:DB:666:A:H2'	23:DB:667:U:C6	2.46	0.50
23:DB:743:A:OP1	26:DD:135:GLY:HA2	2.11	0.50
29:DE:148:ILE:HG22	29:DE:148:ILE:O	2.11	0.50
42:DN:59:SER:O	42:DN:63:ARG:HB2	2.11	0.50
35:DV:80:HIS:HA	35:DV:87:GLN:OE1	2.12	0.50
52:DW:32:ALA:O	52:DW:34:SER:N	2.44	0.50
39:DX:20:ASN:O	39:DX:24:GLU:HB3	2.12	0.50
1:AA:1283:U:O2'	1:AA:1284:C:H5'	2.11	0.50
1:AA:634:C:H2'	1:AA:635:A:H8	1.76	0.50
1:AA:665:A:H2'	1:AA:725:G:N2	2.26	0.50
1:AA:789:U:H2'	1:AA:791:G:OP2	2.11	0.50
1:AA:853:C:C2'	1:AA:854:U:H5'	2.41	0.50
2:AC:194:VAL:HG12	2:AC:195:ILE:N	2.26	0.50
2:AC:185:THR:HG22	2:AC:198:LYS:HA	1.93	0.50
5:AF:84:VAL:HG22	5:AF:85:ILE:H	1.76	0.50
6:AG:74:VAL:HG22	6:AG:143:MET:HG3	1.94	0.50
11:AL:49:ARG:HH12	11:AL:88:ASP:CB	2.24	0.50
14:AQ:58:VAL:HB	14:AQ:74:LEU:CD2	2.41	0.50
23:BB:1174:U:H4'	23:BB:1176:U:N3	2.26	0.50
23:BB:1203:U:H3'	23:BB:1204:A:H5''	1.93	0.50
23:BB:2133:G:H3'	23:BB:2134:A:H5''	1.93	0.50
23:BB:2714:G:O2'	23:BB:2715:C:H5'	2.12	0.50
23:BB:41:C:O2'	23:BB:42:A:H5'	2.10	0.50
23:BB:663:G:OP1	37:BL:17:LYS:HG2	2.11	0.50
23:BB:979:A:H3'	23:BB:980:A:C5'	2.41	0.50
25:BC:129:LEU:HB3	25:BC:134:ILE:HG22	1.94	0.50
25:BC:71:ASP:CA	25:BC:118:GLY:HA2	2.41	0.50
25:BC:90:ILE:HA	25:BC:103:ILE:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:37:VAL:CG2	26:BD:91:THR:HA	2.41	0.50
29:BE:109:LEU:HA	29:BE:112:LEU:HB2	1.92	0.50
47:BF:135:ILE:HG13	47:BF:137:PHE:CD2	2.46	0.50
47:BF:28:PRO:O	47:BF:168:LEU:HD11	2.12	0.50
24:BI:11:GLN:O	24:BI:11:GLN:HG3	2.10	0.50
43:BO:24:THR:O	43:BO:90:VAL:HB	2.11	0.50
28:BP:63:ILE:HG22	28:BP:63:ILE:O	2.10	0.50
50:BT:40:LYS:HA	50:BT:43:ILE:CG2	2.41	0.50
50:BT:67:VAL:HG23	50:BT:75:GLY:O	2.12	0.50
46:BU:86:PHE:HE1	46:BU:88:ASP:HB2	1.75	0.50
1:CA:1280:A:H3'	1:CA:1281:C:H5''	1.93	0.50
1:CA:240:G:OP1	1:CA:240:G:H4'	2.11	0.50
1:CA:204:G:N2	1:CA:466:A:N6	2.60	0.50
1:CA:489:C:H2'	1:CA:490:C:C6	2.46	0.50
1:CA:531:U:H5'	1:CA:531:U:H6	1.76	0.50
1:CA:546:A:P	3:CD:68:GLU:HB3	2.52	0.50
1:CA:765:G:O6	1:CA:811:C:H5	1.93	0.50
2:CC:41:TYR:CZ	2:CC:89:VAL:HG11	2.46	0.50
6:CG:43:TYR:O	6:CG:46:LEU:N	2.44	0.50
8:CI:93:LEU:O	8:CI:97:LEU:HG	2.11	0.50
11:CL:120:ARG:HG2	11:CL:121:PRO:HD2	1.93	0.50
11:CL:31:GLY:HA3	11:CL:54:VAL:HG13	1.94	0.50
36:D2:34:ARG:HH11	36:D2:34:ARG:HG3	1.75	0.50
23:DB:138:U:H2'	23:DB:140:C:O4'	2.12	0.50
23:DB:1516:G:O2'	23:DB:1517:G:H5'	2.11	0.50
23:DB:677:A:H2'	23:DB:678:C:H6	1.75	0.50
23:DB:704:G:H1'	23:DB:727:A:H61	1.77	0.50
25:DC:93:VAL:CG1	25:DC:101:ARG:H	2.24	0.50
25:DC:71:ASP:CA	25:DC:118:GLY:HA2	2.41	0.50
26:DD:51:THR:HG23	26:DD:78:GLY:O	2.11	0.50
47:DF:33:ILE:HG22	47:DF:34:THR:N	2.25	0.50
23:DB:1099:G:H5''	24:DI:2:LYS:HB2	1.93	0.50
41:DJ:17:VAL:CG2	41:DJ:137:PRO:HB2	2.36	0.50
27:DK:47:ILE:HG23	27:DK:49:ARG:H	1.76	0.50
42:DN:3:HIS:O	42:DN:4:ARG:HB2	2.11	0.50
43:DO:51:ALA:CB	43:DO:78:VAL:HG13	2.41	0.50
49:DR:2:TYR:N	49:DR:42:ALA:HB2	2.26	0.50
1:AA:1095:U:H2'	1:AA:1096:C:H6	1.75	0.50
1:AA:853:C:O2'	1:AA:854:U:H5'	2.11	0.50
3:AD:40:HIS:O	3:AD:43:ARG:HG2	2.12	0.50
6:AG:64:ALA:HA	6:AG:67:ASN:HD22	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:5:TYR:HB3	8:AI:88:GLU:OE2	2.11	0.50
9:AJ:6:ILE:CD1	9:AJ:76:ILE:HG13	2.40	0.50
21:AN:28:ALA:HA	21:AN:32:ASP:HB2	1.92	0.50
23:BB:2066:C:O2'	23:BB:2067:G:H5'	2.12	0.50
23:BB:2259:U:H2'	23:BB:2260:C:H6	1.76	0.50
23:BB:2886:A:N6	31:B0:39:ARG:CZ	2.72	0.50
25:BC:108:GLY:C	25:BC:110:LYS:H	2.15	0.50
47:BF:109:ARG:HB3	47:BF:135:ILE:HB	1.94	0.50
47:BF:12:VAL:O	47:BF:16:MET:HG2	2.12	0.50
47:BF:35:LEU:HD23	47:BF:36:ASN:H	1.76	0.50
24:BI:49:GLU:CG	24:BI:54:ILE:HD11	2.41	0.50
23:BB:558:U:P	41:BJ:113:PRO:HG2	2.51	0.50
41:BJ:88:THR:HG22	41:BJ:91:GLU:OE1	2.10	0.50
27:BK:107:LEU:H	27:BK:107:LEU:CD1	2.18	0.50
27:BK:12:ASP:OD2	27:BK:86:LEU:HG	2.11	0.50
28:BP:88:ARG:HB2	28:BP:112:ARG:CZ	2.41	0.50
44:BQ:94:LEU:C	44:BQ:96:ASP:H	2.14	0.50
46:BU:18:LYS:HE2	46:BU:19:GLY:N	2.27	0.50
35:BV:53:LYS:HZ2	35:BV:54:ALA:H	1.57	0.50
52:BW:23:LYS:O	52:BW:66:VAL:HB	2.11	0.50
51:BZ:15:GLY:O	51:BZ:27:ARG:HG3	2.11	0.50
1:CA:1288:A:C2	1:CA:1371:G:H1'	2.45	0.50
1:CA:1400:C:H4'	1:CA:1401:G:OP2	2.10	0.50
1:CA:1434:A:H2'	1:CA:1435:G:O4'	2.11	0.50
1:CA:451:A:H5'	13:CP:70:ARG:HH22	1.74	0.50
1:CA:580:C:H2'	1:CA:581:G:O4'	2.11	0.50
18:CB:20:ARG:HG3	18:CB:20:ARG:HH11	1.77	0.50
3:CD:138:PRO:HA	3:CD:181:PHE:HD2	1.75	0.50
4:CE:132:PRO:O	4:CE:136:VAL:HG12	2.11	0.50
4:CE:150:GLU:CD	4:CE:150:GLU:H	2.15	0.50
5:CF:11:HIS:HB3	5:CF:14:GLN:CG	2.38	0.50
19:CU:29:ALA:HA	19:CU:32:ARG:HG3	1.93	0.50
10:CK:108:ASN:ND2	19:CU:6:ARG:HD2	2.21	0.50
23:DB:1266:G:N2	23:DB:2012:G:H2'	2.26	0.50
23:DB:1458:U:O2'	23:DB:1459:G:H4'	2.11	0.50
23:DB:1508:A:H2'	23:DB:1509:A:C2	2.47	0.50
23:DB:1681:G:H21	23:DB:1762:A:H3'	1.76	0.50
23:DB:18:U:O3'	44:DQ:22:GLY:HA2	2.11	0.50
23:DB:1965:C:H5''	23:DB:1966:A:H2'	1.94	0.50
23:DB:2333:A:H5'	23:DB:2335:A:H1'	1.93	0.50
23:DB:445:C:H2'	23:DB:446:G:O4'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:962:G:H2'	23:DB:963:U:H6	1.76	0.50
23:DB:978:G:O2'	23:DB:979:A:H5'	2.11	0.50
23:DB:2050:C:H1'	26:DD:161:MET:CE	2.41	0.50
29:DE:165:HIS:C	29:DE:167:VAL:H	2.13	0.50
48:DG:118:ALA:C	48:DG:120:ILE:H	2.14	0.50
24:DI:102:ARG:HG3	24:DI:141:ASP:HB2	1.93	0.50
37:DL:81:ASP:HA	37:DL:84:LYS:HE2	1.93	0.50
43:DO:24:THR:O	43:DO:90:VAL:HB	2.11	0.50
50:DT:85:VAL:C	50:DT:86:THR:HG23	2.32	0.50
46:DU:11:ILE:O	46:DU:12:VAL:HB	2.11	0.50
1:AA:978:A:H5'	1:AA:1362:A:N6	2.26	0.50
1:AA:148:G:H2'	1:AA:149:A:O4'	2.11	0.50
1:AA:807:A:H2'	1:AA:808:C:C6	2.47	0.50
18:AB:13:VAL:HG11	18:AB:207:ARG:HH21	1.75	0.50
18:AB:42:LEU:HA	18:AB:45:THR:HB	1.93	0.50
7:AH:48:PHE:HA	7:AH:59:GLU:O	2.12	0.50
11:AL:17:LYS:HD2	11:AL:18:SER:O	2.11	0.50
12:AM:28:ARG:NH2	12:AM:61:LYS:HB2	2.25	0.50
20:AO:81:LEU:HD23	20:AO:82:ILE:N	2.26	0.50
13:AP:41:PRO:O	13:AP:42:ILE:HD13	2.11	0.50
31:B0:38:LEU:HB3	31:B0:41:HIS:NE2	2.26	0.50
23:BB:1930:G:H2'	23:BB:1968:G:N1	2.26	0.50
23:BB:2231:U:H2'	23:BB:2232:C:H6	1.76	0.50
23:BB:234:U:H2'	23:BB:235:U:H6	1.75	0.50
23:BB:2519:U:C6	23:BB:2542:A:N6	2.79	0.50
23:BB:2633:G:H2'	23:BB:2634:A:O4'	2.11	0.50
23:BB:2835:A:H62	23:BB:2878:U:H2'	1.77	0.50
23:BB:441:U:H2'	23:BB:442:G:H8	1.75	0.50
29:BE:28:VAL:O	29:BE:32:VAL:HG22	2.12	0.50
47:BF:101:ARG:NH1	47:BF:138:PRO:HB2	2.26	0.50
47:BF:155:ILE:HG22	47:BF:157:THR:N	2.25	0.50
47:BF:3:LEU:HD11	47:BF:172:PHE:CE1	2.46	0.50
47:BF:91:ARG:O	47:BF:92:GLY:C	2.50	0.50
40:BH:47:PHE:CA	40:BH:50:ARG:HE	2.11	0.50
41:BJ:24:THR:OG1	41:BJ:27:ARG:HD2	2.12	0.50
37:BL:123:ARG:HA	37:BL:143:GLU:CB	2.32	0.50
42:BN:9:GLN:HA	42:BN:17:ARG:HD3	1.92	0.50
49:BR:1:MET:HA	49:BR:42:ALA:HB3	1.92	0.50
44:BQ:108:LEU:HG	49:BR:48:LYS:HG2	1.92	0.50
52:BW:65:LYS:NZ	52:BW:84:GLU:HB3	2.25	0.50
1:CA:1463:U:H2'	1:CA:1464:U:C6	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:187:G:N2	1:CA:189:A:H3'	2.27	0.50
2:CC:184:ASN:O	2:CC:199:VAL:HG22	2.10	0.50
4:CE:104:ILE:HD11	4:CE:111:ARG:HA	1.94	0.50
6:CG:73:GLU:HG2	6:CG:90:VAL:HG12	1.92	0.50
7:CH:77:VAL:HG23	7:CH:126:CYS:HA	1.94	0.50
8:CI:51:LEU:HA	8:CI:54:VAL:CG2	2.33	0.50
21:CN:28:ALA:O	21:CN:32:ASP:HB3	2.12	0.50
14:CQ:58:VAL:HB	14:CQ:74:LEU:CD2	2.41	0.50
23:DB:1039:A:H2'	23:DB:1040:A:H8	1.75	0.50
23:DB:1567:G:H1'	23:DB:1568:G:C6	2.47	0.50
23:DB:1271:G:OP2	23:DB:1648:U:OP1	2.28	0.50
23:DB:1689:A:H2'	23:DB:1690:A:C8	2.47	0.50
23:DB:1838:C:H4'	23:DB:1839:G:H8	1.76	0.50
23:DB:2221:G:H2'	23:DB:2222:C:H6	1.75	0.50
23:DB:2789:C:H2'	23:DB:2893:A:N7	2.27	0.50
23:DB:547:A:N3	23:DB:547:A:C2'	2.66	0.50
23:DB:547:A:H5'	23:DB:548:G:C4	2.47	0.50
23:DB:8:C:O2'	23:DB:9:G:H5'	2.10	0.50
26:DD:117:GLY:HA2	26:DD:164:GLN:OE1	2.11	0.50
26:DD:202:ILE:HG22	26:DD:204:LYS:HZ2	1.75	0.50
29:DE:105:LEU:HD21	29:DE:177:PRO:HA	1.92	0.50
47:DF:62:GLN:NE2	47:DF:90:LEU:HA	2.25	0.50
47:DF:64:PRO:HA	47:DF:88:VAL:CG1	2.41	0.50
40:DH:94:ILE:HG23	40:DH:98:ASP:CB	2.42	0.50
24:DI:75:ALA:O	24:DI:79:LEU:HG	2.12	0.50
27:DK:107:LEU:CD1	27:DK:107:LEU:H	2.21	0.50
38:DM:111:GLU:HA	38:DM:114:ARG:HH22	1.76	0.50
45:DS:45:VAL:HG23	45:DS:46:LEU:N	2.26	0.50
52:DW:17:ALA:HA	52:DW:35:ILE:CG2	2.31	0.50
1:AA:1139:G:O4'	1:AA:1140:C:H5	1.95	0.50
1:AA:1182:G:H4'	1:AA:1183:U:C5'	2.40	0.50
1:AA:1381:U:O2	6:AG:78:ARG:HB2	2.12	0.50
1:AA:1499:A:H2'	1:AA:1500:A:C8	2.39	0.50
1:AA:1512:U:H2'	1:AA:1513:A:C8	2.46	0.50
1:AA:161:A:H2'	1:AA:162:A:C8	2.47	0.50
1:AA:56:U:H2'	1:AA:57:G:C8	2.45	0.50
18:AB:71:THR:HG23	18:AB:92:ASN:O	2.10	0.50
2:AC:127:VAL:HG23	2:AC:128:MET:H	1.76	0.50
3:AD:152:SER:CA	3:AD:155:LYS:HD3	2.36	0.50
3:AD:2:ARG:H	3:AD:4:LEU:HD11	1.77	0.50
8:AI:94:ARG:CB	8:AI:94:ARG:HH11	2.23	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AN:61:ASN:HB3	21:AN:72:PHE:CZ	2.47	0.50
14:AQ:25:GLU:N	14:AQ:40:THR:HG23	2.27	0.50
23:BB:1083:U:H2'	23:BB:1085:A:OP2	2.11	0.50
23:BB:1170:C:O5'	23:BB:1170:C:H6	1.95	0.50
23:BB:1629:U:O2	23:BB:2698:U:H5''	2.11	0.50
23:BB:1854:A:H62	23:BB:1888:G:H1'	1.75	0.50
23:BB:2040:G:H2'	23:BB:2041:U:O4'	2.11	0.50
23:BB:2221:G:H2'	23:BB:2222:C:H6	1.76	0.50
23:BB:2249:U:N3	23:BB:2253:G:OP2	2.41	0.50
23:BB:2339:C:H2'	23:BB:2340:A:C8	2.47	0.50
23:BB:2754:U:H6	23:BB:2754:U:O5'	1.94	0.50
23:BB:2787:C:H1'	26:BD:63:PRO:HG3	1.94	0.50
23:BB:2839:G:O2'	23:BB:2840:C:H5'	2.11	0.50
23:BB:2888:C:H2'	23:BB:2889:C:C6	2.47	0.50
23:BB:471:A:H2'	23:BB:472:A:O4'	2.12	0.50
23:BB:57:C:H2'	23:BB:58:G:H8	1.76	0.50
23:BB:850:U:H2'	23:BB:851:C:H6	1.75	0.50
47:BF:127:TYR:O	47:BF:155:ILE:HB	2.12	0.50
22:BA:43:C:O2'	47:BF:91:ARG:HD2	2.11	0.50
24:BI:23:VAL:HG23	24:BI:24:GLY:H	1.76	0.50
24:BI:33:ASN:HD21	24:BI:64:ARG:NH1	2.03	0.50
38:BM:57:VAL:O	38:BM:57:VAL:HG12	2.11	0.50
42:BN:101:GLY:CA	42:BN:110:MET:H	2.15	0.50
42:BN:65:LEU:O	42:BN:68:ALA:HB3	2.12	0.50
28:BP:6:GLN:NE2	28:BP:7:LEU:HD12	2.26	0.50
45:BS:18:ARG:HB3	45:BS:76:VAL:CG2	2.42	0.50
52:BW:59:PHE:O	52:BW:60:ALA:CB	2.59	0.50
23:BB:1158:C:H4'	30:BY:30:ARG:HH21	1.77	0.50
1:CA:1157:A:H5'	1:CA:1158:C:C6	2.46	0.50
1:CA:236:A:O2'	1:CA:237:G:H5'	2.11	0.50
1:CA:251:G:H22	1:CA:272:C:N4	2.09	0.50
1:CA:377:G:H2'	1:CA:378:G:C8	2.45	0.50
1:CA:89:U:H2'	1:CA:90:C:C6	2.47	0.50
18:CB:131:LYS:HG3	18:CB:132:GLU:N	2.27	0.50
18:CB:14:HIS:O	18:CB:15:PHE:C	2.50	0.50
18:CB:17:HIS:HB3	18:CB:187:ASP:OD2	2.12	0.50
5:CF:53:LYS:NZ	5:CF:53:LYS:H	2.08	0.50
6:CG:57:GLU:O	6:CG:61:PHE:N	2.44	0.50
7:CH:10:LEU:HA	7:CH:74:ILE:HD11	1.94	0.50
13:CP:74:LEU:O	13:CP:78:VAL:HG12	2.12	0.50
22:DA:49:C:P	43:DO:30:ARG:HH12	2.34	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1007:C:H4'	41:DJ:110:PRO:HB3	1.92	0.50
23:DB:1113:U:H2'	23:DB:1114:C:C6	2.47	0.50
23:DB:1341:G:H5'	50:DT:61:LEU:HD13	1.94	0.50
23:DB:1790:C:H2'	23:DB:1791:A:C8	2.47	0.50
23:DB:1886:U:O5'	23:DB:1886:U:H6	1.95	0.50
23:DB:2052:A:OP1	26:DD:145:SER:HA	2.11	0.50
23:DB:2356:U:H5''	52:DW:16:GLU:HG3	1.93	0.50
23:DB:2822:G:H2'	23:DB:2823:A:H5''	1.93	0.50
23:DB:2888:C:H2'	23:DB:2889:C:C6	2.47	0.50
23:DB:569:U:H2'	23:DB:570:G:O4'	2.12	0.50
23:DB:599:A:O2'	23:DB:600:G:H5'	2.12	0.50
25:DC:108:GLY:C	25:DC:110:LYS:H	2.15	0.50
25:DC:12:ARG:HA	25:DC:15:VAL:CG2	2.42	0.50
26:DD:61:THR:OG1	26:DD:64:GLU:HB2	2.11	0.50
47:DF:32:LYS:H	47:DF:95:MET:CE	2.25	0.50
23:DB:2009:A:O4'	42:DN:107:ASN:HB3	2.11	0.50
50:DT:28:ASN:ND2	50:DT:29:THR:HG23	2.27	0.50
50:DT:40:LYS:O	50:DT:43:ILE:HG22	2.11	0.50
23:DB:922:C:H1'	52:DW:22:VAL:HG21	1.94	0.50
1:AA:1073:U:H4'	18:AB:104:LYS:HE2	1.93	0.50
1:AA:1111:A:C2	2:AC:176:THR:HG23	2.47	0.50
1:AA:1338:G:H2'	1:AA:1339:A:H8	1.75	0.50
1:AA:1389:C:H2'	1:AA:1390:U:C6	2.46	0.50
1:AA:404:G:O2'	1:AA:405:U:H5'	2.12	0.50
1:AA:531:U:C6	1:AA:531:U:H5'	2.47	0.50
1:AA:562:U:H4'	1:AA:563:A:O5'	2.11	0.50
1:AA:891:U:O2'	1:AA:892:A:H5'	2.11	0.50
18:AB:131:LYS:HB3	18:AB:131:LYS:HZ2	1.76	0.50
2:AC:72:PRO:HB3	2:AC:102:ILE:HG21	1.94	0.50
3:AD:90:LEU:HD21	3:AD:196:GLU:CB	2.41	0.50
5:AF:3:HIS:CG	5:AF:92:THR:HG23	2.46	0.50
5:AF:92:THR:HG22	5:AF:94:HIS:N	2.16	0.50
8:AI:27:ILE:HG23	8:AI:64:ILE:HD11	1.92	0.50
8:AI:50:PRO:HG2	8:AI:51:LEU:H	1.76	0.50
10:AK:53:GLY:O	10:AK:56:LYS:HB3	2.11	0.50
12:AM:28:ARG:NH2	12:AM:62:PHE:HB2	2.24	0.50
12:AM:38:ILE:HG13	12:AM:55:LEU:CD2	2.38	0.50
12:AM:3:ILE:HD12	12:AM:7:ASN:O	2.12	0.50
16:AS:48:ILE:HB	16:AS:59:VAL:HG23	1.92	0.50
17:AT:49:ALA:HA	17:AT:52:GLU:HB3	1.92	0.50
34:B3:14:LYS:HG2	34:B3:14:LYS:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:B4:1:MET:CE	32:B4:36:ARG:HB2	2.42	0.50
23:BB:1866:A:H2'	23:BB:1867:G:O4'	2.12	0.50
23:BB:1902:C:H2'	23:BB:1903:G:O4'	2.12	0.50
23:BB:1946:U:H2'	23:BB:1947:C:C6	2.46	0.50
23:BB:2466:C:O2'	23:BB:2467:C:H5'	2.11	0.50
23:BB:195:A:H1'	23:BB:250:G:N2	2.27	0.50
23:BB:974:G:P	49:BR:78:ARG:HD3	2.52	0.50
29:BE:58:LYS:HZ3	29:BE:58:LYS:H	1.55	0.50
47:BF:32:LYS:H	47:BF:95:MET:CE	2.25	0.50
48:BG:106:LEU:C	48:BG:108:PHE:H	2.14	0.50
27:BK:47:ILE:HG23	27:BK:49:ARG:H	1.76	0.50
37:BL:30:THR:HB	37:BL:34:GLY:O	2.10	0.50
28:BP:111:GLU:OE1	28:BP:113:LEU:HB2	2.10	0.50
28:BP:63:ILE:HG12	28:BP:68:GLY:HA2	1.94	0.50
23:BB:18:U:O3'	44:BQ:22:GLY:HA2	2.11	0.50
44:BQ:30:VAL:HG22	44:BQ:31:TYR:N	2.23	0.50
45:BS:66:ILE:CA	45:BS:69:LEU:HD13	2.41	0.50
46:BU:5:ARG:HH22	46:BU:93:ARG:HD3	1.76	0.50
1:CA:356:A:H1'	1:CA:368:U:O2'	2.12	0.50
1:CA:601:G:H2'	1:CA:602:A:C8	2.46	0.50
1:CA:630:A:H2'	1:CA:631:C:C6	2.46	0.50
1:CA:645:G:H2'	1:CA:646:G:H8	1.76	0.50
18:CB:50:ASN:HA	18:CB:53:LEU:HB3	1.92	0.50
18:CB:83:ALA:O	18:CB:88:GLN:HB2	2.10	0.50
2:CC:106:ARG:O	2:CC:107:LYS:C	2.49	0.50
12:CM:78:ARG:O	12:CM:82:LEU:HG	2.11	0.50
16:CS:11:ASP:O	16:CS:14:LEU:HG	2.12	0.50
19:CU:16:ARG:HH12	19:CU:19:LYS:HZ3	1.59	0.50
22:DA:100:G:H2'	22:DA:101:A:O4'	2.11	0.50
23:DB:1047:G:O3'	23:DB:1048:A:H8	1.94	0.50
23:DB:126:A:H5'	36:D2:19:ARG:CG	2.40	0.50
23:DB:1312:U:H4'	23:DB:1313:U:O5'	2.12	0.50
23:DB:1403:A:O2'	23:DB:1404:C:H5'	2.11	0.50
23:DB:2290:G:H2'	23:DB:2291:U:C6	2.47	0.50
23:DB:2751:G:H3'	23:DB:2752:C:C6	2.47	0.50
23:DB:572:A:H5''	49:DR:80:ARG:NH2	2.27	0.50
23:DB:803:U:O2'	23:DB:804:A:H5'	2.12	0.50
23:DB:90:U:H2'	23:DB:91:A:C2	2.46	0.50
47:DF:133:GLU:HA	47:DF:150:GLY:CA	2.41	0.50
41:DJ:58:ASN:HA	41:DJ:127:GLY:CA	2.42	0.50
44:DQ:30:VAL:HG22	44:DQ:31:TYR:N	2.23	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DQ:42:GLY:HA3	49:DR:75:VAL:CG2	2.41	0.50
50:DT:31:VAL:HA	50:DT:83:ALA:HB3	1.93	0.50
50:DT:7:LEU:O	50:DT:7:LEU:HD13	2.12	0.50
46:DU:33:VAL:HG13	46:DU:66:VAL:HB	1.93	0.50
52:DW:65:LYS:NZ	52:DW:84:GLU:HB3	2.26	0.50
30:DY:58:GLU:H	30:DY:58:GLU:CD	2.15	0.50
1:AA:1100:C:C2	1:AA:1102:A:H5'	2.47	0.50
1:AA:1215:G:H5'	1:AA:1215:G:C8	2.45	0.50
1:AA:1392:G:H2'	1:AA:1393:U:C6	2.46	0.50
1:AA:1499:A:OP2	1:AA:1499:A:H3'	2.11	0.50
1:AA:68:G:C5'	1:AA:171:A:H1'	2.40	0.50
1:AA:683:G:O2'	1:AA:684:U:H5'	2.10	0.50
2:AC:68:HIS:HA	2:AC:103:ALA:HB3	1.94	0.50
6:AG:110:ARG:HB3	6:AG:112:ASP:OD1	2.12	0.50
6:AG:149:ALA:HB2	10:AK:60:PHE:HB2	1.93	0.50
6:AG:3:ARG:NH1	6:AG:3:ARG:HB3	2.27	0.50
9:AJ:35:GLN:HG2	9:AJ:78:GLU:OE1	2.12	0.50
21:AN:77:GLY:O	21:AN:78:LEU:HD12	2.12	0.50
13:AP:38:PHE:O	13:AP:50:THR:HG23	2.11	0.50
17:AT:73:ARG:HG3	17:AT:74:HIS:N	2.25	0.50
19:AU:29:ALA:HA	19:AU:32:ARG:HG3	1.94	0.50
23:BB:1000:A:H2'	23:BB:1001:A:H8	1.76	0.50
23:BB:1080:A:H2'	23:BB:1081:U:H6	1.77	0.50
23:BB:1403:A:O2'	23:BB:1404:C:H5'	2.12	0.50
23:BB:1818:U:C4	25:BC:152:GLN:HB3	2.47	0.50
23:BB:2102:G:H2'	23:BB:2103:C:C6	2.47	0.50
23:BB:2544:G:O2'	23:BB:2545:G:H5'	2.12	0.50
23:BB:6:A:H4'	41:BJ:131:ASN:O	2.12	0.50
23:BB:708:G:H2'	23:BB:709:U:C6	2.46	0.50
23:BB:970:U:H2'	23:BB:971:G:H8	1.76	0.50
23:BB:1657:U:O2'	26:BD:138:LEU:HD12	2.12	0.50
23:BB:37:C:O2'	29:BE:45:ALA:HA	2.12	0.50
47:BF:117:SER:HB3	47:BF:120:SER:HB3	1.93	0.50
47:BF:148:VAL:O	47:BF:149:ARG:HG2	2.11	0.50
47:BF:30:VAL:HG21	47:BF:96:TRP:HE1	1.76	0.50
48:BG:51:PHE:CD2	48:BG:68:ARG:HG2	2.47	0.50
40:BH:53:GLU:C	40:BH:54:LEU:HD23	2.32	0.50
41:BJ:104:ALA:O	41:BJ:108:MET:HG3	2.12	0.50
37:BL:89:VAL:O	37:BL:89:VAL:HG13	2.12	0.50
45:BS:48:LYS:O	45:BS:51:LEU:HB3	2.11	0.50
50:BT:31:VAL:HA	50:BT:83:ALA:HB3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BX:20:ASN:HD22	39:BX:20:ASN:N	2.09	0.50
1:CA:1175:G:O2'	1:CA:1176:A:H5'	2.12	0.50
1:CA:130:A:H1'	1:CA:263:A:O2'	2.11	0.50
1:CA:14:U:H2'	1:CA:14:U:O2	2.11	0.50
1:CA:1511:G:O2'	1:CA:1512:U:H5'	2.12	0.50
1:CA:413:G:O6	3:CD:32:LYS:HE2	2.11	0.50
1:CA:489:C:H2'	1:CA:490:C:H6	1.77	0.50
1:CA:547:A:H4'	1:CA:548:G:O5'	2.11	0.50
1:CA:756:C:H2'	1:CA:757:U:O4'	2.12	0.50
2:CC:128:MET:H	2:CC:128:MET:CE	2.24	0.50
2:CC:26:LYS:HG3	2:CC:27:GLU:HG3	1.94	0.50
2:CC:52:SER:OG	2:CC:53:ARG:N	2.42	0.50
2:CC:18:ASN:C	2:CC:55:VAL:HG13	2.32	0.50
4:CE:63:MET:O	4:CE:66:ALA:HB3	2.11	0.50
6:CG:119:LEU:O	6:CG:123:LEU:HG	2.11	0.50
6:CG:24:LYS:O	6:CG:28:ILE:HG12	2.12	0.50
10:CK:30:ILE:HG22	10:CK:45:THR:CB	2.42	0.50
10:CK:22:ILE:HG21	10:CK:95:THR:HG21	1.92	0.50
16:CS:38:THR:HG22	16:CS:39:ILE:O	2.12	0.50
22:DA:52:A:O2'	22:DA:53:A:H5'	2.11	0.50
23:DB:1636:U:H2'	23:DB:1637:A:H8	1.76	0.50
23:DB:1664:A:H1'	23:DB:2726:A:N1	2.27	0.50
23:DB:170:U:H2'	23:DB:171:U:H6	1.77	0.50
23:DB:2250:G:N2	38:DM:82:MET:HB3	2.27	0.50
23:DB:2369:A:H2'	23:DB:2370:G:C8	2.47	0.50
23:DB:2645:G:H4'	23:DB:2732:G:H2'	1.93	0.50
23:DB:2772:C:H4'	26:DD:171:THR:CG2	2.41	0.50
47:DF:147:ARG:O	47:DF:148:VAL:C	2.50	0.50
24:DI:76:ALA:HA	24:DI:135:MET:SD	2.51	0.50
24:DI:21:PRO:CB	24:DI:22:PRO:HD3	2.38	0.50
41:DJ:25:LEU:O	41:DJ:27:ARG:N	2.43	0.50
27:DK:71:ARG:CB	27:DK:72:PRO:HD2	2.37	0.50
37:DL:30:THR:HB	37:DL:34:GLY:O	2.12	0.50
38:DM:126:ILE:N	38:DM:126:ILE:HD12	2.25	0.50
49:DR:34:GLU:OE1	49:DR:60:LYS:HG2	2.11	0.50
1:AA:1105:A:O2'	1:AA:1106:G:H5'	2.12	0.50
1:AA:149:A:H1'	1:AA:1446:A:C2	2.46	0.50
1:AA:216:U:H2'	1:AA:217:C:C6	2.46	0.50
1:AA:356:A:H1'	1:AA:368:U:O2'	2.12	0.50
1:AA:737:C:H2'	1:AA:738:C:H6	1.76	0.50
1:AA:747:A:C4	1:AA:748:G:H1'	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:122:GLN:OE1	2:AC:127:VAL:HG21	2.12	0.50
2:AC:149:LYS:CB	2:AC:200:TRP:HB2	2.35	0.50
2:AC:38:VAL:HG23	2:AC:39:ARG:N	2.27	0.50
7:AH:43:GLY:HA2	7:AH:63:LYS:NZ	2.27	0.50
11:AL:49:ARG:HH12	11:AL:88:ASP:HB3	1.77	0.50
1:AA:135:C:O2	13:AP:1:MET:HB2	2.11	0.50
36:B2:3:ARG:HE	36:B2:4:THR:H	1.60	0.50
23:BB:1237:A:H2'	23:BB:1237:A:N3	2.27	0.50
23:BB:1255:U:C6	29:BE:68:ALA:HB2	2.47	0.50
23:BB:1550:C:H2'	23:BB:1551:A:H8	1.77	0.50
23:BB:1655:A:H2'	23:BB:1656:C:O4'	2.12	0.50
23:BB:1666:G:O3'	27:BK:6:THR:HG23	2.11	0.50
23:BB:2259:U:H2'	23:BB:2260:C:C6	2.47	0.50
23:BB:2849:U:N3	23:BB:2867:G:C8	2.80	0.50
23:BB:2896:C:H2'	23:BB:2897:U:C6	2.47	0.50
23:BB:95:A:H4'	39:BX:38:GLN:O	2.12	0.50
25:BC:77:VAL:HG22	25:BC:113:ASP:O	2.11	0.50
29:BE:32:VAL:HG23	29:BE:33:VAL:N	2.26	0.50
48:BG:148:ARG:CD	48:BG:152:ARG:HH11	2.22	0.50
48:BG:17:LYS:HG3	48:BG:19:ASN:OD1	2.12	0.50
48:BG:9:VAL:HA	48:BG:48:THR:CB	2.39	0.50
40:BH:116:ARG:HH21	40:BH:139:PHE:HB3	1.76	0.50
38:BM:66:ARG:CZ	38:BM:101:VAL:HG11	2.42	0.50
42:BN:103:ARG:HB3	42:BN:106:ASP:O	2.12	0.50
42:BN:92:GLY:O	42:BN:94:TYR:N	2.44	0.50
35:BV:29:ILE:HD13	35:BV:31:TYR:CE2	2.47	0.50
52:BW:19:ARG:HD3	52:BW:36:ILE:HD11	1.93	0.50
39:BX:31:GLN:O	39:BX:37:LEU:HB2	2.11	0.50
30:BY:43:ILE:HG13	30:BY:44:ARG:N	2.26	0.50
1:CA:1020:G:C2'	1:CA:1021:A:H5'	2.42	0.50
1:CA:1049:U:H1'	1:CA:1201:A:C5	2.46	0.50
1:CA:174:A:O2'	1:CA:175:C:H5'	2.12	0.50
1:CA:382:A:H2'	1:CA:383:A:H8	1.77	0.50
1:CA:408:A:H3'	1:CA:409:U:C6	2.47	0.50
1:CA:430:A:O2'	1:CA:431:A:H5'	2.12	0.50
3:CD:146:GLU:CD	3:CD:146:GLU:N	2.65	0.50
1:CA:8:A:N6	3:CD:53:GLN:HE22	2.10	0.50
4:CE:113:VAL:CG2	4:CE:114:LEU:H	2.18	0.50
4:CE:81:GLN:H	4:CE:146:MET:CE	2.17	0.50
5:CF:35:LYS:O	5:CF:64:VAL:HG13	2.12	0.50
6:CG:119:LEU:HG	6:CG:123:LEU:HD11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:CN:56:PRO:O	21:CN:58:ARG:N	2.45	0.50
16:CS:38:THR:HA	16:CS:68:HIS:O	2.12	0.50
22:DA:106:G:H2'	22:DA:107:G:O4'	2.11	0.50
22:DA:117:G:O2'	22:DA:118:C:H5'	2.11	0.50
23:DB:1098:A:P	24:DI:3:LYS:HG2	2.52	0.50
23:DB:1535:A:H2	23:DB:1538:G:H1'	1.76	0.50
23:DB:1561:C:H2'	23:DB:1562:U:C6	2.47	0.50
23:DB:1597:A:C5'	23:DB:1598:A:H5'	2.34	0.50
23:DB:182:A:O2'	23:DB:183:C:H5'	2.11	0.50
23:DB:1949:G:H2'	23:DB:1950:G:C8	2.47	0.50
23:DB:2040:G:H2'	23:DB:2041:U:O4'	2.10	0.50
23:DB:2081:U:H2'	23:DB:2082:A:C8	2.47	0.50
23:DB:231:A:H3'	23:DB:232:G:H8	1.75	0.50
23:DB:2345:G:H5'	23:DB:2347:C:O4'	2.12	0.50
23:DB:2352:A:C6	52:DW:30:VAL:HG11	2.47	0.50
23:DB:278:A:O3'	23:DB:279:A:H8	1.94	0.50
23:DB:335:C:O2'	23:DB:336:C:H5'	2.12	0.50
23:DB:831:G:O2'	23:DB:832:U:H5'	2.12	0.50
25:DC:189:ALA:O	25:DC:190:THR:C	2.49	0.50
29:DE:60:TRP:CZ3	29:DE:69:ARG:HA	2.47	0.50
47:DF:109:ARG:HG3	47:DF:137:PHE:CB	2.41	0.50
49:DR:83:TYR:HE2	49:DR:85:LYS:HE3	1.75	0.50
50:DT:69:ARG:HE	50:DT:69:ARG:CA	2.22	0.50
50:DT:69:ARG:HG3	50:DT:70:HIS:N	2.26	0.50
46:DU:18:LYS:HE2	46:DU:19:GLY:N	2.27	0.50
1:AA:270:A:H2'	1:AA:271:C:H6	1.73	0.50
1:AA:279:A:H5'	1:AA:281:G:C5'	2.41	0.50
1:AA:470:C:H2'	1:AA:471:U:H6	1.76	0.50
1:AA:8:A:H5"	4:AE:125:LYS:HB3	1.94	0.50
6:AG:74:VAL:HA	6:AG:87:PRO:HA	1.94	0.50
7:AH:124:ILE:C	7:AH:125:ILE:HD12	2.31	0.50
9:AJ:22:THR:OG1	9:AJ:72:ARG:HG3	2.11	0.50
31:B0:53:VAL:O	31:B0:54:ILE:HB	2.12	0.50
23:BB:116:C:H2'	23:BB:117:G:O4'	2.11	0.50
23:BB:1516:G:O2'	23:BB:1517:G:H5'	2.11	0.50
23:BB:1774:C:C2'	23:BB:1774:C:O2	2.59	0.50
23:BB:1911:U:H2'	23:BB:1918:A:C2	2.46	0.50
23:BB:917:A:H5"	23:BB:2268:A:N6	2.27	0.50
26:BD:32:ASN:HD22	26:BD:50:VAL:HG21	1.76	0.50
23:BB:797:G:OP2	29:BE:57:LYS:HB2	2.12	0.50
47:BF:147:ARG:O	47:BF:148:VAL:C	2.50	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:68:ARG:HB2	40:BH:134:VAL:CG1	2.41	0.50
40:BH:99:ILE:HD11	40:BH:128:HIS:CG	2.47	0.50
42:BN:19:ALA:C	42:BN:21:PHE:N	2.65	0.50
46:BU:62:ALA:O	46:BU:63:ALA:HB3	2.12	0.50
52:BW:33:GLY:O	52:BW:34:SER:HB2	2.12	0.50
51:BZ:5:CYS:HB3	51:BZ:9:GLY:H	1.77	0.50
1:CA:1291:U:O2'	1:CA:1292:G:H5'	2.10	0.50
1:CA:292:G:O2'	1:CA:609:A:N6	2.45	0.50
1:CA:462:G:H5'	1:CA:463:U:P	2.52	0.50
1:CA:57:G:H2'	1:CA:58:C:H6	1.76	0.50
1:CA:691:G:H1'	1:CA:696:A:H61	1.77	0.50
1:CA:731:G:O2'	1:CA:732:C:H5'	2.12	0.50
1:CA:939:G:H4'	6:CG:101:ARG:NH2	2.26	0.50
18:CB:32:GLY:O	18:CB:39:ILE:HB	2.12	0.50
2:CC:161:ILE:N	2:CC:161:ILE:HD12	2.26	0.50
8:CI:15:ALA:O	8:CI:67:LYS:HD2	2.12	0.50
9:CJ:57:VAL:O	9:CJ:58:ASN:HB2	2.12	0.50
9:CJ:81:GLU:N	9:CJ:82:LYS:HZ3	2.10	0.50
20:CO:71:LYS:NZ	20:CO:72:ARG:N	2.60	0.50
13:CP:38:PHE:O	13:CP:50:THR:HG23	2.12	0.50
17:CT:24:ARG:CG	17:CT:65:LEU:HD11	2.42	0.50
19:CU:40:PRO:C	19:CU:42:THR:H	2.15	0.50
34:D3:57:VAL:C	34:D3:59:ALA:H	2.15	0.50
23:DB:1099:G:H4'	24:DI:4:VAL:HG12	1.93	0.50
23:DB:1175:A:H2'	23:DB:1176:U:O4'	2.12	0.50
23:DB:1324:G:H1'	23:DB:1616:A:C6	2.47	0.50
23:DB:1365:A:N3	23:DB:1365:A:H2'	2.27	0.50
23:DB:1731:G:O2'	23:DB:1732:C:H5''	2.11	0.50
23:DB:1998:A:O2'	23:DB:1999:C:H5'	2.12	0.50
23:DB:2104:C:H2'	23:DB:2105:U:H6	1.77	0.50
23:DB:2144:G:O2'	23:DB:2146:C:H5''	2.12	0.50
23:DB:2297:A:C2	23:DB:2320:U:H4'	2.47	0.50
23:DB:2415:G:H2'	23:DB:2416:C:H6	1.77	0.50
23:DB:2521:C:H2'	23:DB:2522:U:C6	2.47	0.50
23:DB:2676:C:H2'	23:DB:2677:G:C8	2.47	0.50
23:DB:496:G:C1'	45:DS:61:ASN:HD21	2.25	0.50
23:DB:920:A:H2'	23:DB:921:C:C6	2.47	0.50
26:DD:117:GLY:HA2	26:DD:164:GLN:CD	2.32	0.50
37:DL:125:LEU:HB2	37:DL:143:GLU:OE2	2.12	0.50
37:DL:19:LEU:HD22	37:DL:31:GLY:O	2.12	0.50
23:DB:2723:C:H4'	42:DN:1:MET:SD	2.51	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DW:59:PHE:O	52:DW:60:ALA:CB	2.60	0.50
52:DW:69:GLU:HG3	52:DW:80:SER:OG	2.11	0.50
1:AA:131:A:H2'	1:AA:132:C:C6	2.46	0.49
1:AA:1375:A:H2'	1:AA:1376:U:O4'	2.11	0.49
1:AA:1413:A:O2'	1:AA:1414:U:H5'	2.12	0.49
1:AA:575:G:H4'	1:AA:576:C:O5'	2.12	0.49
1:AA:795:C:H5''	10:AK:127:ARG:HH21	1.77	0.49
1:AA:932:C:H4'	6:AG:3:ARG:NH2	2.26	0.49
18:AB:19:THR:HG23	18:AB:20:ARG:N	2.26	0.49
2:AC:102:ILE:N	2:AC:102:ILE:HD12	2.27	0.49
3:AD:197:HIS:O	3:AD:201:GLU:HG3	2.11	0.49
5:AF:68:GLN:HG2	5:AF:69:GLU:N	2.26	0.49
6:AG:110:ARG:HB2	6:AG:118:ARG:HB3	1.94	0.49
7:AH:25:THR:O	7:AH:26:MET:HB3	2.11	0.49
8:AI:40:ARG:H	8:AI:44:ARG:HD3	1.76	0.49
10:AK:30:ILE:HG22	10:AK:45:THR:CB	2.42	0.49
12:AM:21:ILE:CG2	12:AM:23:GLY:H	2.22	0.49
21:AN:43:ALA:HA	21:AN:46:LYS:CG	2.42	0.49
14:AQ:32:ILE:HG23	14:AQ:33:TYR:CD2	2.46	0.49
16:AS:28:LYS:N	16:AS:28:LYS:HD2	2.27	0.49
1:AA:1314:C:H41	16:AS:3:SER:HB3	1.77	0.49
34:B3:50:SER:C	34:B3:52:GLY:H	2.15	0.49
22:BA:60:C:H2'	22:BA:61:G:C8	2.44	0.49
23:BB:1886:U:H6	23:BB:1886:U:O5'	1.95	0.49
23:BB:2394:C:OP1	37:BL:63:LYS:HG2	2.12	0.49
23:BB:78:U:H2'	23:BB:79:C:C6	2.47	0.49
23:BB:836:G:H2'	23:BB:837:C:C6	2.47	0.49
25:BC:184:GLU:O	25:BC:186:ASP:N	2.40	0.49
25:BC:144:GLU:OE2	25:BC:188:ARG:HG3	2.11	0.49
29:BE:138:LEU:HD22	29:BE:143:LEU:HB2	1.94	0.49
29:BE:145:ASP:H	29:BE:166:LYS:CB	2.24	0.49
48:BG:148:ARG:HA	48:BG:161:VAL:HB	1.93	0.49
48:BG:83:THR:C	48:BG:84:LYS:HD3	2.32	0.49
37:BL:14:LYS:O	37:BL:16:GLY:N	2.45	0.49
43:BO:51:ALA:CB	43:BO:78:VAL:HG13	2.42	0.49
50:BT:50:LEU:C	50:BT:52:GLU:H	2.15	0.49
46:BU:13:LEU:HA	46:BU:18:LYS:CE	2.40	0.49
1:CA:103:U:H2'	1:CA:104:G:O4'	2.11	0.49
1:CA:1386:G:O3'	54:CA:1659:SCM:H1M2	2.12	0.49
1:CA:1395:C:O2'	1:CA:1396:A:H5'	2.12	0.49
1:CA:1499:A:OP2	1:CA:1499:A:H3'	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:24:U:O2'	1:CA:25:C:H5'	2.12	0.49
1:CA:891:U:O2'	1:CA:892:A:H5'	2.11	0.49
1:CA:911:U:H2'	1:CA:912:C:C6	2.47	0.49
2:CC:145:ALA:C	2:CC:147:GLY:H	2.15	0.49
3:CD:123:MET:SD	3:CD:127:ARG:N	2.85	0.49
3:CD:25:ARG:O	3:CD:27:ILE:HG13	2.12	0.49
5:CF:53:LYS:HZ2	5:CF:53:LYS:H	1.60	0.49
9:CJ:7:ARG:HG3	9:CJ:102:LEU:C	2.32	0.49
10:CK:36:ARG:HG3	10:CK:36:ARG:HH11	1.77	0.49
21:CN:86:ALA:HA	21:CN:89:ARG:HH21	1.75	0.49
16:CS:18:VAL:HG13	16:CS:19:GLU:N	2.26	0.49
16:CS:59:VAL:C	16:CS:60:PHE:HD2	2.14	0.49
22:DA:116:G:H4'	43:DO:54:VAL:HG22	1.94	0.49
23:DB:1045:C:C3'	23:DB:1046:A:H5'	2.42	0.49
23:DB:1441:G:H2'	23:DB:1442:U:H6	1.73	0.49
23:DB:1809:A:H2'	23:DB:1810:A:C8	2.47	0.49
23:DB:1878:G:H2'	23:DB:1879:C:C6	2.47	0.49
23:DB:37:C:H4'	23:DB:451:U:OP1	2.12	0.49
23:DB:710:U:H2'	23:DB:711:G:C8	2.47	0.49
25:DC:34:GLU:O	25:DC:34:GLU:HG3	2.12	0.49
26:DD:149:ASN:C	26:DD:152:PRO:HD2	2.31	0.49
26:DD:11:MET:H	26:DD:25:THR:HA	1.77	0.49
29:DE:126:VAL:HG22	29:DE:127:GLU:N	2.20	0.49
29:DE:189:THR:O	29:DE:193:VAL:HG23	2.11	0.49
47:DF:100:GLU:O	47:DF:104:THR:HB	2.12	0.49
48:DG:152:ARG:HD3	48:DG:153:PRO:CD	2.40	0.49
28:DP:111:GLU:OE1	28:DP:113:LEU:HB2	2.12	0.49
45:DS:36:LEU:HD11	45:DS:47:VAL:HB	1.94	0.49
30:DY:35:VAL:HG11	30:DY:37:ARG:HH12	1.77	0.49
1:AA:1113:C:O2'	1:AA:1114:C:H5'	2.11	0.49
1:AA:123:U:H5''	1:AA:311:C:O2'	2.13	0.49
1:AA:1340:A:H2'	1:AA:1341:U:O4'	2.12	0.49
1:AA:542:G:O2'	1:AA:543:U:H5'	2.12	0.49
1:AA:962:C:H2'	1:AA:963:G:H8	1.76	0.49
18:AB:57:ASN:HB3	18:AB:219:THR:CG2	2.39	0.49
2:AC:116:ALA:O	2:AC:120:THR:HG23	2.12	0.49
2:AC:153:SER:OG	2:AC:195:ILE:HG23	2.12	0.49
2:AC:13:ILE:O	2:AC:15:LYS:N	2.38	0.49
8:AI:29:ILE:HG23	8:AI:64:ILE:HG21	1.94	0.49
11:AL:85:ARG:CB	11:AL:93:ARG:HA	2.42	0.49
12:AM:33:LEU:HD13	12:AM:39:ALA:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AN:27:LYS:HA	21:AN:31:SER:HB2	1.92	0.49
17:AT:19:HIS:O	17:AT:23:ARG:HG2	2.11	0.49
36:B2:9:VAL:CG1	36:B2:10:LEU:N	2.75	0.49
23:BB:1026:G:H2'	23:BB:1027:A:C8	2.48	0.49
23:BB:1203:U:O4'	37:BL:3:LEU:HD12	2.12	0.49
23:BB:2720:U:H5''	28:BP:52:ARG:HH22	1.74	0.49
23:BB:281:C:H2'	23:BB:282:A:C8	2.47	0.49
23:BB:1812:U:H1'	25:BC:43:ASN:HD21	1.77	0.49
47:BF:33:ILE:HG22	47:BF:34:THR:N	2.27	0.49
48:BG:122:ALA:CA	48:BG:132:LEU:HA	2.41	0.49
41:BJ:15:TRP:HB3	41:BJ:137:PRO:HB3	1.93	0.49
41:BJ:1:MET:HG2	41:BJ:2:LYS:HG2	1.94	0.49
49:BR:74:ILE:HB	49:BR:87:GLN:O	2.11	0.49
52:BW:9:THR:C	52:BW:10:ARG:HD3	2.32	0.49
52:BW:81:ILE:O	52:BW:81:ILE:HG13	2.12	0.49
1:CA:1366:C:H2'	1:CA:1367:C:C6	2.47	0.49
1:CA:580:C:H2'	1:CA:581:G:C8	2.46	0.49
1:CA:706:A:H4'	10:CK:30:ILE:CD1	2.42	0.49
1:CA:665:A:H2'	1:CA:725:G:N2	2.27	0.49
1:CA:767:A:H2'	1:CA:768:A:H8	1.77	0.49
1:CA:83:C:OP1	1:CA:83:C:H4'	2.12	0.49
18:CB:124:THR:O	18:CB:127:LYS:HE3	2.11	0.49
18:CB:68:PHE:HA	18:CB:161:PHE:O	2.12	0.49
18:CB:185:ILE:HA	18:CB:199:ILE:O	2.12	0.49
18:CB:64:GLY:O	18:CB:66:ILE:N	2.45	0.49
5:CF:3:HIS:CD2	5:CF:3:HIS:N	2.80	0.49
7:CH:45:ILE:C	7:CH:63:LYS:HE3	2.32	0.49
1:CA:568:G:O6	11:CL:1:ALA:HB2	2.12	0.49
21:CN:40:ARG:CZ	16:CS:6:LYS:HB2	2.41	0.49
31:D0:53:VAL:O	31:D0:54:ILE:HB	2.12	0.49
36:D2:3:ARG:HE	36:D2:4:THR:H	1.60	0.49
23:DB:123:G:H2'	23:DB:124:G:O4'	2.12	0.49
23:DB:1306:C:O2'	23:DB:1307:A:H5'	2.13	0.49
23:DB:1454:C:C5	42:DN:64:ARG:HG2	2.47	0.49
23:DB:1636:U:H2'	23:DB:1637:A:C8	2.47	0.49
23:DB:1728:C:H2'	23:DB:1730:C:O2	2.11	0.49
23:DB:189:G:H2'	23:DB:205:G:N2	2.26	0.49
23:DB:2357:G:H22	23:DB:2359:C:H3'	1.77	0.49
23:DB:2443:C:H2'	23:DB:2444:G:C8	2.47	0.49
23:DB:2462:C:H2'	23:DB:2463:C:H6	1.77	0.49
23:DB:2636:C:H4'	26:DD:81:GLU:OE2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2639:A:H2'	23:DB:2640:G:O4'	2.11	0.49
23:DB:736:C:H2'	23:DB:737:C:H6	1.76	0.49
23:DB:757:G:H2'	23:DB:758:C:H5'	1.93	0.49
26:DD:133:THR:HG23	26:DD:134:HIS:CD2	2.47	0.49
26:DD:202:ILE:HD12	26:DD:202:ILE:N	2.27	0.49
29:DE:47:LYS:HB3	29:DE:51:GLU:HB2	1.93	0.49
47:DF:37:MET:HA	47:DF:151:LEU:HA	1.94	0.49
47:DF:91:ARG:O	47:DF:92:GLY:C	2.50	0.49
40:DH:148:ALA:O	40:DH:149:GLU:HG2	2.12	0.49
41:DJ:24:THR:OG1	41:DJ:27:ARG:HD2	2.12	0.49
27:DK:12:ASP:OD2	27:DK:86:LEU:HG	2.12	0.49
38:DM:10:ARG:HG3	38:DM:10:ARG:HH21	1.77	0.49
38:DM:41:LEU:HB3	38:DM:46:ILE:HG21	1.94	0.49
42:DN:47:VAL:O	42:DN:50:PRO:HD2	2.12	0.49
42:DN:58:ASP:O	42:DN:59:SER:HB3	2.11	0.49
42:DN:65:LEU:O	42:DN:68:ALA:HB3	2.13	0.49
45:DS:13:SER:OG	45:DS:14:ALA:N	2.45	0.49
46:DU:35:VAL:CG2	46:DU:38:ILE:HG21	2.41	0.49
46:DU:49:PRO:HA	46:DU:53:GLN:HE21	1.77	0.49
52:DW:18:LYS:HA	52:DW:36:ILE:HG12	1.93	0.49
39:DX:39:GLN:HB3	39:DX:41:HIS:NE2	2.27	0.49
39:DX:39:GLN:HB3	39:DX:42:LEU:HD13	1.94	0.49
1:AA:1176:A:H2'	1:AA:1177:G:O4'	2.11	0.49
1:AA:590:U:H2'	1:AA:591:U:H6	1.78	0.49
1:AA:691:G:H1'	1:AA:696:A:H61	1.77	0.49
1:AA:687:A:C2	1:AA:704:A:C5	3.00	0.49
1:AA:706:A:H4'	10:AK:30:ILE:HD11	1.94	0.49
3:AD:54:LEU:HD11	3:AD:55:ARG:HH21	1.78	0.49
6:AG:46:LEU:O	6:AG:57:GLU:HB3	2.13	0.49
8:AI:48:ARG:O	8:AI:52:GLU:HB2	2.13	0.49
1:AA:1124:G:C5'	9:AJ:38:GLY:HA3	2.42	0.49
1:AA:706:A:H4'	10:AK:30:ILE:CD1	2.41	0.49
11:AL:32:VAL:HB	11:AL:55:ARG:HB3	1.94	0.49
20:AO:43:PHE:CE2	20:AO:53:ARG:HA	2.47	0.49
13:AP:5:ARG:HH21	13:AP:24:SER:HA	1.73	0.49
15:AR:32:ILE:HD11	15:AR:58:ILE:HG23	1.95	0.49
36:B2:34:ARG:HG3	36:B2:34:ARG:HH11	1.76	0.49
23:BB:1526:C:H2'	23:BB:1527:G:O4'	2.12	0.49
23:BB:1973:G:O2'	23:BB:1974:C:H5'	2.13	0.49
23:BB:210:C:O2'	23:BB:211:C:H5'	2.13	0.49
23:BB:2369:A:H2'	23:BB:2370:G:C8	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2547:A:H4'	27:BK:29:HIS:NE2	2.27	0.49
23:BB:2694:G:H2'	23:BB:2695:U:C6	2.48	0.49
23:BB:414:C:H2'	23:BB:415:A:H8	1.76	0.49
23:BB:579:G:H4'	23:BB:2018:G:H5''	1.95	0.49
25:BC:140:VAL:HG12	25:BC:141:HIS:N	2.23	0.49
25:BC:145:MET:HG3	25:BC:152:GLN:OE1	2.12	0.49
25:BC:41:GLY:HA3	25:BC:53:ILE:CG2	2.37	0.49
29:BE:136:GLN:HB2	29:BE:139:LYS:HE3	1.94	0.49
47:BF:116:LEU:HD23	47:BF:174:PHE:CE2	2.48	0.49
48:BG:102:ILE:O	48:BG:113:ASP:HA	2.12	0.49
48:BG:167:VAL:HG23	48:BG:168:VAL:H	1.77	0.49
23:BB:1060:U:C4	24:BI:131:THR:HG22	2.47	0.49
44:BQ:59:LEU:C	44:BQ:59:LEU:HD13	2.32	0.49
45:BS:76:VAL:HG12	45:BS:103:ILE:HA	1.94	0.49
1:CA:1435:G:H2'	1:CA:1436:U:C6	2.47	0.49
1:CA:501:C:H1'	1:CA:549:C:H1'	1.94	0.49
1:CA:531:U:H5'	1:CA:531:U:C6	2.48	0.49
1:CA:590:U:H2'	1:CA:591:U:C6	2.47	0.49
1:CA:721:G:H4'	1:CA:722:G:O4'	2.12	0.49
3:CD:138:PRO:HA	3:CD:181:PHE:CD2	2.47	0.49
4:CE:149:PRO:HG2	4:CE:150:GLU:H	1.78	0.49
6:CG:112:ASP:H	6:CG:118:ARG:NE	2.10	0.49
8:CI:70:GLY:O	8:CI:74:GLN:HB2	2.12	0.49
11:CL:106:VAL:HG23	11:CL:116:TYR:HB3	1.95	0.49
11:CL:86:VAL:CG2	11:CL:89:LEU:HB2	2.41	0.49
14:CQ:18:LYS:HE3	14:CQ:48:GLU:HG2	1.93	0.49
23:DB:1370:C:H2'	23:DB:1371:G:O4'	2.12	0.49
23:DB:1386:C:H2'	23:DB:1387:A:C8	2.48	0.49
23:DB:1390:U:O2'	23:DB:1391:U:H5'	2.12	0.49
23:DB:699:A:H4'	23:DB:1634:A:N7	2.27	0.49
23:DB:1819:A:OP1	25:DC:159:THR:HG21	2.13	0.49
23:DB:1854:A:H62	23:DB:1888:G:H1'	1.76	0.49
23:DB:2054:A:H2'	31:D0:4:GLN:OE1	2.11	0.49
23:DB:2056:G:H21	31:D0:1:ALA:HA	1.76	0.49
23:DB:2075:U:H2'	23:DB:2238:G:N2	2.28	0.49
23:DB:2259:U:H2'	23:DB:2260:C:H6	1.78	0.49
23:DB:230:G:H2'	23:DB:231:A:H8	1.77	0.49
23:DB:2740:A:H2'	23:DB:2741:A:C8	2.47	0.49
23:DB:340:A:H2'	23:DB:341:C:C5'	2.42	0.49
23:DB:531:C:O2'	23:DB:563:A:H5''	2.13	0.49
47:DF:7:TYR:O	47:DF:12:VAL:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:147:ARG:O	47:DF:147:ARG:HD2	2.12	0.49
47:DF:87:LYS:C	47:DF:88:VAL:HG23	2.32	0.49
24:DI:85:ILE:HD12	24:DI:87:SER:O	2.12	0.49
27:DK:17:ARG:HB3	27:DK:45:GLU:HG3	1.95	0.49
27:DK:35:VAL:CG2	27:DK:36:GLY:H	2.13	0.49
38:DM:64:TRP:HB2	38:DM:104:GLU:HB2	1.94	0.49
42:DN:28:LEU:CA	42:DN:34:ILE:HD11	2.42	0.49
23:DB:2820:A:OP1	42:DN:4:ARG:HA	2.12	0.49
43:DO:76:LYS:HG3	43:DO:113:ALA:HB2	1.94	0.49
45:DS:26:GLY:O	45:DS:28:LYS:N	2.45	0.49
1:AA:1077:G:H2'	1:AA:1079:G:OP2	2.12	0.49
1:AA:14:U:H2'	1:AA:14:U:O2	2.13	0.49
1:AA:253:A:H2'	1:AA:254:G:H8	1.76	0.49
1:AA:398:U:H2'	1:AA:399:G:H8	1.76	0.49
1:AA:420:U:O2'	1:AA:421:U:H5''	2.13	0.49
18:AB:67:LEU:HA	18:AB:89:PHE:O	2.11	0.49
2:AC:76:ILE:HA	2:AC:83:VAL:CG2	2.36	0.49
4:AE:156:ARG:HB2	7:AH:43:GLY:HA3	1.93	0.49
34:B3:7:ARG:HG3	34:B3:7:ARG:NH1	2.27	0.49
22:BA:49:C:P	43:BO:30:ARG:HH12	2.36	0.49
23:BB:1128:G:O2'	23:BB:1129:A:H5''	2.12	0.49
23:BB:1190:G:H5''	37:BL:32:GLY:O	2.12	0.49
23:BB:1569:A:H2'	23:BB:1570:A:C8	2.47	0.49
23:BB:1811:G:O2'	23:BB:1812:U:H5'	2.12	0.49
23:BB:2256:G:H2'	23:BB:2257:U:H6	1.77	0.49
23:BB:2547:A:H2'	23:BB:2548:U:H6	1.76	0.49
23:BB:256:A:H2'	23:BB:257:C:H6	1.78	0.49
23:BB:877:A:H5''	23:BB:877:A:H8	1.77	0.49
26:BD:125:TRP:CD2	26:BD:160:LYS:HB3	2.48	0.49
26:BD:202:ILE:HG22	26:BD:204:LYS:NZ	2.26	0.49
47:BF:139:GLU:CG	47:BF:140:ILE:H	2.25	0.49
47:BF:57:ALA:HB2	47:BF:64:PRO:CG	2.42	0.49
37:BL:3:LEU:O	37:BL:5:THR:N	2.44	0.49
38:BM:126:ILE:HD12	38:BM:126:ILE:N	2.27	0.49
50:BT:40:LYS:O	50:BT:43:ILE:HG22	2.13	0.49
46:BU:6:ARG:O	46:BU:24:VAL:HB	2.13	0.49
1:CA:1121:U:O2'	1:CA:1122:U:H5'	2.13	0.49
1:CA:1167:A:H2'	1:CA:1169:A:N6	2.27	0.49
1:CA:218:U:H2'	1:CA:219:U:C6	2.46	0.49
1:CA:241:G:O2'	1:CA:242:G:H5'	2.12	0.49
1:CA:403:C:O2'	1:CA:404:G:H5'	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:502:A:H2'	1:CA:503:C:C6	2.47	0.49
1:CA:505:G:H4'	1:CA:534:U:C4	2.47	0.49
1:CA:692:U:H2'	1:CA:694:A:OP2	2.12	0.49
1:CA:687:A:C2	1:CA:704:A:C5	3.00	0.49
3:CD:41:GLY:C	3:CD:43:ARG:H	2.16	0.49
9:CJ:8:ILE:CD1	9:CJ:76:ILE:HG23	2.43	0.49
10:CK:14:GLN:OE1	10:CK:77:GLY:HA3	2.13	0.49
12:CM:64:VAL:HA	12:CM:68:LEU:HD12	1.94	0.49
14:CQ:24:ILE:N	14:CQ:24:ILE:HD12	2.27	0.49
16:CS:55:GLN:HA	16:CS:55:GLN:NE2	2.26	0.49
36:D2:10:LEU:HD21	36:D2:14:ARG:CZ	2.42	0.49
34:D3:14:LYS:HG2	34:D3:14:LYS:O	2.12	0.49
23:DB:1518:C:H2'	23:DB:1519:G:C8	2.48	0.49
23:DB:151:C:H2'	23:DB:152:A:C8	2.48	0.49
23:DB:1661:G:O2'	23:DB:1662:U:H5'	2.12	0.49
23:DB:2221:G:O2'	23:DB:2222:C:H5'	2.12	0.49
23:DB:252:G:O2'	23:DB:253:C:H5'	2.12	0.49
23:DB:38:A:O2'	29:DE:43:THR:HA	2.13	0.49
23:DB:702:U:H2'	23:DB:703:U:C6	2.47	0.49
23:DB:899:A:H2'	23:DB:900:A:O4'	2.12	0.49
25:DC:91:ALA:CB	25:DC:105:ALA:HB2	2.42	0.49
26:DD:138:LEU:N	26:DD:138:LEU:HD22	2.28	0.49
29:DE:4:VAL:HG12	29:DE:5:LEU:N	2.22	0.49
24:DI:92:PRO:O	24:DI:93:ASN:HB2	2.12	0.49
41:DJ:114:LEU:O	41:DJ:118:MET:HG3	2.13	0.49
42:DN:72:ASP:C	42:DN:74:GLU:H	2.16	0.49
26:DD:15:PHE:CE2	28:DP:77:SER:HB2	2.48	0.49
46:DU:48:VAL:O	46:DU:48:VAL:HG13	2.12	0.49
35:DV:59:GLU:HG2	35:DV:60:VAL:N	2.28	0.49
52:DW:23:LYS:HD2	52:DW:24:ARG:N	2.25	0.49
1:AA:1048:G:O3'	1:AA:1049:U:H3'	2.12	0.49
1:AA:1124:G:H5''	9:AJ:37:ARG:O	2.13	0.49
1:AA:1275:A:H2'	1:AA:1276:G:O4'	2.12	0.49
1:AA:1400:C:H4'	1:AA:1401:G:OP2	2.13	0.49
1:AA:1435:G:H2'	1:AA:1436:U:C6	2.47	0.49
1:AA:34:C:H2'	1:AA:35:G:C8	2.47	0.49
1:AA:394:G:O2'	1:AA:395:C:H5'	2.12	0.49
1:AA:903:G:H2'	1:AA:904:U:C6	2.48	0.49
1:AA:996:A:H2'	1:AA:997:U:H6	1.75	0.49
2:AC:134:LYS:HE3	2:AC:167:TYR:OH	2.12	0.49
2:AC:94:ALA:O	2:AC:96:VAL:HG22	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:170:LEU:HD12	3:AD:170:LEU:O	2.13	0.49
5:AF:6:ILE:HG13	5:AF:62:MET:HB2	1.94	0.49
7:AH:101:ALA:HB3	7:AH:112:ASP:HB3	1.95	0.49
7:AH:99:GLY:HA2	7:AH:129:ALA:HA	1.95	0.49
1:AA:1250:A:H4'	8:AI:69:GLY:O	2.12	0.49
9:AJ:73:LEU:HD13	9:AJ:75:ASP:HB2	1.95	0.49
12:AM:63:VAL:O	12:AM:68:LEU:HD21	2.11	0.49
21:AN:63:CYS:HB2	21:AN:79:SER:HB3	1.93	0.49
10:AK:110:THR:CG2	19:AU:4:LYS:HA	2.42	0.49
34:B3:21:PHE:HB2	34:B3:49:VAL:HG22	1.93	0.49
22:BA:112:G:O2'	22:BA:113:C:H5'	2.13	0.49
23:BB:1449:G:O2'	23:BB:1450:G:H5'	2.13	0.49
23:BB:1508:A:H2'	23:BB:1509:A:C2	2.47	0.49
23:BB:1731:G:O2'	23:BB:1732:C:H5''	2.12	0.49
23:BB:1842:G:H2'	23:BB:1843:C:C6	2.48	0.49
23:BB:189:G:H2'	23:BB:205:G:H22	1.77	0.49
23:BB:2653:U:H3'	23:BB:2654:A:H2'	1.95	0.49
23:BB:2789:C:H2'	23:BB:2893:A:N7	2.27	0.49
23:BB:596:U:H2'	23:BB:597:G:H8	1.78	0.49
23:BB:693:A:H2'	23:BB:694:U:H6	1.78	0.49
23:BB:753:A:H2'	23:BB:754:U:H6	1.73	0.49
25:BC:173:LEU:H	25:BC:173:LEU:HD22	1.77	0.49
25:BC:202:ARG:HE	25:BC:213:ARG:NH2	2.09	0.49
47:BF:111:ARG:HH11	47:BF:135:ILE:HG21	1.77	0.49
47:BF:29:ARG:H	47:BF:29:ARG:HE	1.59	0.49
40:BH:47:PHE:HA	40:BH:50:ARG:NE	2.09	0.49
24:BI:89:SER:HA	24:BI:97:VAL:CG2	2.42	0.49
41:BJ:124:VAL:O	41:BJ:125:TYR:HB2	2.13	0.49
37:BL:136:GLU:C	37:BL:138:ALA:H	2.15	0.49
38:BM:100:LYS:CD	38:BM:101:VAL:H	2.26	0.49
43:BO:52:SER:C	43:BO:54:VAL:H	2.15	0.49
28:BP:25:VAL:HB	28:BP:84:SER:O	2.13	0.49
27:BK:64:ARG:HH21	28:BP:67:GLU:HG3	1.78	0.49
49:BR:38:VAL:HG22	49:BR:40:MET:H	1.78	0.49
50:BT:15:HIS:N	50:BT:32:LEU:HA	2.17	0.49
46:BU:78:LYS:CD	46:BU:79:ALA:H	2.24	0.49
1:CA:1118:U:H1'	1:CA:1179:A:C4	2.47	0.49
1:CA:1203:C:H4'	21:CN:66:THR:HG22	1.95	0.49
1:CA:121:U:H3'	1:CA:121:U:OP1	2.12	0.49
1:CA:1312:G:H2'	1:CA:1313:U:C6	2.47	0.49
1:CA:1374:A:O2'	1:CA:1375:A:H5'	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1426:G:O2'	1:CA:1427:C:H5'	2.13	0.49
1:CA:1476:A:H2'	1:CA:1477:U:O4'	2.12	0.49
1:CA:1486:G:H2'	1:CA:1487:G:O4'	2.12	0.49
1:CA:597:G:H2'	1:CA:598:U:H5'	1.94	0.49
1:CA:607:A:H2'	1:CA:608:A:C8	2.46	0.49
1:CA:861:G:H2'	1:CA:862:C:C6	2.46	0.49
18:CB:71:THR:C	18:CB:76:SER:HB3	2.32	0.49
2:CC:104:GLU:HG2	2:CC:105:VAL:N	2.27	0.49
3:CD:2:ARG:H	3:CD:4:LEU:HD11	1.78	0.49
7:CH:17:GLN:HE21	7:CH:62:LEU:HG	1.78	0.49
1:CA:1347:G:H2'	8:CI:109:GLN:O	2.13	0.49
8:CI:44:ARG:HB2	8:CI:48:ARG:CZ	2.43	0.49
8:CI:90:ASP:O	8:CI:92:SER:N	2.43	0.49
14:CQ:45:VAL:HG12	14:CQ:46:HIS:N	2.27	0.49
16:CS:14:LEU:HD22	16:CS:34:SER:HB2	1.94	0.49
16:CS:47:THR:HA	16:CS:60:PHE:CZ	2.47	0.49
32:D4:1:MET:CE	32:D4:36:ARG:HB2	2.42	0.49
23:DB:1459:G:H2'	23:DB:1461:C:C4	2.47	0.49
23:DB:1494:A:H2'	23:DB:1495:A:C8	2.48	0.49
23:DB:783:A:H4'	23:DB:1779:U:O2	2.11	0.49
23:DB:2047:C:H2'	23:DB:2048:G:C8	2.47	0.49
23:DB:2350:C:H2'	23:DB:2351:G:C8	2.48	0.49
23:DB:11:C:N4	23:DB:2629:U:H3	2.10	0.49
23:DB:2677:G:H2'	23:DB:2678:C:H6	1.74	0.49
23:DB:2831:G:H1'	23:DB:2883:A:C2	2.47	0.49
23:DB:475:C:H4'	23:DB:509:C:H2'	1.94	0.49
23:DB:574:A:H4'	23:DB:575:A:C5'	2.43	0.49
25:DC:201:LEU:O	25:DC:202:ARG:O	2.29	0.49
29:DE:105:LEU:HD21	29:DE:177:PRO:HB3	1.94	0.49
29:DE:32:VAL:HG23	29:DE:33:VAL:N	2.27	0.49
47:DF:155:ILE:HG22	47:DF:156:THR:N	2.27	0.49
47:DF:78:ILE:HA	47:DF:82:TYR:CD1	2.47	0.49
40:DH:47:PHE:O	40:DH:51:ARG:HG3	2.13	0.49
40:DH:5:LEU:HD13	40:DH:12:LEU:C	2.33	0.49
23:DB:7:G:H5'	41:DJ:132:HIS:CD2	2.48	0.49
27:DK:73:ASP:OD2	27:DK:75:SER:HB3	2.12	0.49
37:DL:78:ARG:NH2	37:DL:113:ALA:HB1	2.27	0.49
38:DM:54:THR:C	38:DM:56:ALA:H	2.15	0.49
28:DP:26:GLU:HG3	28:DP:43:GLU:HB2	1.93	0.49
28:DP:63:ILE:HA	28:DP:67:GLU:O	2.12	0.49
45:DS:24:ILE:CD1	45:DS:36:LEU:HD21	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DU:3:LYS:CB	46:DU:82:VAL:HG21	2.37	0.49
46:DU:5:ARG:HH22	46:DU:93:ARG:HD3	1.76	0.49
35:DV:63:ILE:HB	35:DV:70:ILE:HG12	1.94	0.49
52:DW:35:ILE:O	52:DW:35:ILE:HG12	2.11	0.49
52:DW:43:LYS:HD2	52:DW:79:ILE:CD1	2.33	0.49
1:AA:1053:G:C6	1:AA:1200:C:H5'	2.48	0.49
1:AA:922:G:N3	1:AA:1398:A:H2	2.10	0.49
1:AA:130:A:H1'	1:AA:263:A:O2'	2.12	0.49
1:AA:721:G:H4'	1:AA:722:G:O4'	2.12	0.49
1:AA:908:A:O2'	1:AA:909:A:H5'	2.12	0.49
18:AB:95:TRP:CZ3	18:AB:171:ALA:HA	2.47	0.49
5:AF:66:ALA:HB1	5:AF:70:VAL:HG21	1.94	0.49
6:AG:4:ARG:NH1	6:AG:5:VAL:H	2.09	0.49
8:AI:90:ASP:HB3	8:AI:93:LEU:HG	1.93	0.49
10:AK:20:ALA:O	10:AK:22:ILE:HG13	2.12	0.49
10:AK:30:ILE:HG22	10:AK:45:THR:HB	1.94	0.49
21:AN:26:LEU:O	21:AN:30:ILE:N	2.43	0.49
14:AQ:45:VAL:HG12	14:AQ:46:HIS:N	2.27	0.49
17:AT:47:GLN:HE21	17:AT:82:ILE:CD1	2.25	0.49
22:BA:16:G:O2'	22:BA:17:C:H5'	2.13	0.49
22:BA:2:G:H2'	22:BA:3:C:C5	2.48	0.49
23:BB:1631:G:H22	23:BB:1633:G:H3'	1.76	0.49
23:BB:2081:U:H2'	23:BB:2082:A:C8	2.47	0.49
23:BB:222:A:N6	23:BB:232:G:H1'	2.27	0.49
23:BB:2599:G:H2'	23:BB:2600:A:H8	1.77	0.49
23:BB:2800:A:H2'	23:BB:2801:G:C4'	2.42	0.49
23:BB:340:A:H2'	23:BB:341:C:C5'	2.43	0.49
23:BB:425:G:O2'	23:BB:426:C:H5'	2.13	0.49
23:BB:692:C:H2'	23:BB:693:A:H8	1.76	0.49
25:BC:244:VAL:HG12	25:BC:250:GLN:N	2.28	0.49
47:BF:134:GLN:HB3	47:BF:149:ARG:CB	2.43	0.49
24:BI:72:THR:HG21	24:BI:111:THR:O	2.13	0.49
28:BP:99:LEU:HD13	28:BP:99:LEU:O	2.12	0.49
44:BQ:59:LEU:HD13	44:BQ:60:TRP:N	2.28	0.49
30:BY:2:LYS:HE2	30:BY:39:ASP:OD2	2.12	0.49
1:CA:1037:C:O2'	1:CA:1038:C:H5'	2.13	0.49
1:CA:1118:U:H2'	1:CA:1119:C:H6	1.78	0.49
1:CA:1254:A:O4'	1:CA:1356:G:H5'	2.13	0.49
1:CA:313:A:H2'	1:CA:314:C:H6	1.78	0.49
1:CA:555:U:H2'	1:CA:556:C:H6	1.78	0.49
1:CA:585:G:O2'	1:CA:586:C:H5'	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:594:U:H2'	1:CA:595:A:O4'	2.13	0.49
1:CA:68:G:H2'	1:CA:69:G:O4'	2.13	0.49
1:CA:859:G:O2'	1:CA:860:A:H5'	2.13	0.49
18:CB:51:GLU:HG2	18:CB:197:PHE:HD1	1.78	0.49
2:CC:155:ARG:H	2:CC:162:ALA:CA	2.13	0.49
2:CC:18:ASN:CA	2:CC:55:VAL:HG22	2.38	0.49
3:CD:72:ARG:HA	3:CD:203:TYR:HE1	1.77	0.49
9:CJ:86:ALA:O	9:CJ:90:LEU:HG	2.13	0.49
11:CL:30:ARG:O	11:CL:57:THR:HG23	2.13	0.49
13:CP:5:ARG:HH21	13:CP:24:SER:HA	1.75	0.49
23:DB:2886:A:N6	31:D0:39:ARG:CZ	2.76	0.49
23:DB:2884:U:C5	31:D0:49:ARG:HG2	2.47	0.49
23:DB:1163:G:O2'	23:DB:1164:C:H5'	2.12	0.49
23:DB:1259:G:H2'	23:DB:1260:A:C8	2.47	0.49
23:DB:138:U:O2'	23:DB:140:C:H5'	2.13	0.49
23:DB:142:A:H2	50:DT:2:ILE:HG22	1.76	0.49
23:DB:1742:U:O2'	23:DB:1743:G:H5'	2.12	0.49
23:DB:1838:C:H4'	23:DB:1839:G:C8	2.47	0.49
23:DB:2394:C:H2'	23:DB:2395:C:C6	2.48	0.49
23:DB:2432:A:O2'	23:DB:2433:A:H5'	2.12	0.49
23:DB:2881:U:O3'	42:DN:96:ARG:HG3	2.13	0.49
23:DB:5:A:H2'	23:DB:6:A:H8	1.72	0.49
23:DB:876:C:H2'	23:DB:876:C:O2	2.11	0.49
25:DC:9:SER:C	25:DC:11:GLY:H	2.16	0.49
47:DF:107:VAL:N	47:DF:108:PRO:CD	2.75	0.49
40:DH:115:VAL:HA	40:DH:132:PHE:HA	1.93	0.49
24:DI:24:GLY:HA2	24:DI:34:ILE:HD12	1.95	0.49
27:DK:108:ARG:HG3	27:DK:108:ARG:O	2.12	0.49
37:DL:18:ARG:O	37:DL:19:LEU:HD12	2.12	0.49
34:D3:23:HIS:CG	37:DL:61:LEU:HD23	2.48	0.49
38:DM:66:ARG:CZ	38:DM:101:VAL:HG11	2.43	0.49
31:D0:53:VAL:CB	42:DN:118:ARG:HH22	2.25	0.49
44:DQ:105:PHE:HA	44:DQ:108:LEU:CD1	2.34	0.49
44:DQ:59:LEU:C	44:DQ:59:LEU:HD13	2.33	0.49
45:DS:88:ARG:HG3	45:DS:88:ARG:HH21	1.77	0.49
46:DU:98:ASN:OD1	46:DU:100:GLU:HB2	2.13	0.49
46:DU:48:VAL:O	46:DU:50:ALA:N	2.45	0.49
35:DV:75:GLN:HB3	35:DV:90:ASP:HB2	1.93	0.49
1:AA:1005:A:H2'	1:AA:1006:G:O4'	2.12	0.49
1:AA:1005:A:H5'	1:AA:1037:C:O2	2.13	0.49
1:AA:1060:U:C5'	9:AJ:53:ILE:HG12	2.40	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1135:U:H5'	1:AA:1136:C:OP1	2.12	0.49
1:AA:1330:U:H5''	12:AM:22:TYR:O	2.12	0.49
1:AA:408:A:H3'	1:AA:409:U:C6	2.45	0.49
1:AA:505:G:H4'	1:AA:534:U:C4	2.46	0.49
1:AA:692:U:O2	1:AA:694:A:H5''	2.13	0.49
1:AA:746:A:H2'	1:AA:747:A:H8	1.76	0.49
1:AA:814:A:O2'	1:AA:815:A:H3'	2.12	0.49
18:AB:180:ILE:O	18:AB:182:VAL:HG23	2.13	0.49
18:AB:22:TRP:HA	18:AB:189:ASN:CA	2.42	0.49
6:AG:109:LYS:HE3	6:AG:109:LYS:O	2.12	0.49
6:AG:29:LEU:HD11	6:AG:119:LEU:HD21	1.94	0.49
10:AK:19:VAL:N	10:AK:34:THR:O	2.46	0.49
10:AK:75:GLU:N	10:AK:75:GLU:CD	2.66	0.49
11:AL:120:ARG:HG2	11:AL:121:PRO:HD2	1.95	0.49
11:AL:21:PRO:C	11:AL:23:LEU:H	2.16	0.49
12:AM:10:ASP:HB3	12:AM:45:SER:HB3	1.94	0.49
1:AA:1227:A:H1'	12:AM:113:LYS:HZ3	1.77	0.49
12:AM:89:ARG:HH12	12:AM:101:THR:CG2	2.26	0.49
21:AN:10:VAL:HB	21:AN:11:LYS:NZ	2.27	0.49
13:AP:28:ARG:HD3	13:AP:29:ASN:ND2	2.28	0.49
17:AT:27:MET:O	17:AT:31:ILE:HG13	2.12	0.49
17:AT:75:LYS:O	17:AT:78:LEU:HB2	2.13	0.49
31:B0:27:LEU:HD13	31:B0:37:HIS:O	2.13	0.49
36:B2:18:PHE:CE1	36:B2:22:MET:HG3	2.48	0.49
23:BB:1054:A:H2'	23:BB:1055:G:O4'	2.11	0.49
23:BB:1157:G:O2'	30:BY:31:ILE:HD13	2.13	0.49
23:BB:1205:A:N7	29:BE:165:HIS:CG	2.81	0.49
23:BB:1327:A:H2'	23:BB:1328:A:H5'	1.94	0.49
23:BB:1392:A:H2'	23:BB:1393:A:C8	2.47	0.49
23:BB:1652:A:H62	42:BN:11:ASN:ND2	2.02	0.49
23:BB:1654:A:H61	23:BB:2049:G:P	2.36	0.49
23:BB:1654:A:H2'	23:BB:1655:A:H8	1.77	0.49
23:BB:1667:G:OP1	27:BK:6:THR:HA	2.12	0.49
23:BB:171:U:H2'	23:BB:172:A:H8	1.73	0.49
23:BB:2100:G:C6	23:BB:2190:G:C6	3.00	0.49
23:BB:2639:A:H2'	23:BB:2640:G:O4'	2.12	0.49
23:BB:2722:G:H2'	23:BB:2723:C:C6	2.47	0.49
23:BB:548:G:H5''	23:BB:549:G:H5'	1.93	0.49
23:BB:921:C:H2'	23:BB:922:C:H6	1.78	0.49
25:BC:79:ARG:HD2	25:BC:81:GLU:OE2	2.12	0.49
29:BE:62:GLN:CG	29:BE:63:LYS:H	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BG:97:VAL:CG2	48:BG:124:CYS:HB2	2.42	0.49
48:BG:16:VAL:HG11	48:BG:44:HIS:NE2	2.28	0.49
40:BH:111:ALA:O	40:BH:132:PHE:HE1	1.94	0.49
40:BH:5:LEU:CD1	40:BH:17:ASP:HB3	2.38	0.49
37:BL:93:ASN:HD22	37:BL:94:THR:H	1.59	0.49
31:B0:53:VAL:CB	42:BN:118:ARG:HH22	2.25	0.49
42:BN:79:LEU:O	42:BN:80:PHE:HB2	2.11	0.49
28:BP:4:ILE:O	28:BP:6:GLN:N	2.40	0.49
45:BS:81:SER:HB3	45:BS:99:ARG:HB3	1.94	0.49
1:CA:1285:A:HO2'	1:CA:1286:U:P	2.35	0.49
1:CA:1355:G:H2'	1:CA:1356:G:H8	1.77	0.49
1:CA:315:A:H5''	1:CA:317:U:OP2	2.12	0.49
1:CA:35:G:H2'	1:CA:36:C:H6	1.78	0.49
1:CA:429:U:C1'	1:CA:430:A:H5''	2.43	0.49
1:CA:44:A:O2'	1:CA:45:G:H5'	2.12	0.49
1:CA:590:U:H2'	1:CA:591:U:H6	1.77	0.49
1:CA:664:G:N2	1:CA:741:G:H1	2.03	0.49
1:CA:853:C:C2'	1:CA:854:U:H5'	2.42	0.49
2:CC:111:ASP:O	2:CC:112:ALA:C	2.50	0.49
3:CD:110:ARG:HG3	3:CD:110:ARG:HH11	1.78	0.49
4:CE:92:ARG:O	4:CE:126:ALA:HB1	2.12	0.49
6:CG:51:GLN:O	6:CG:51:GLN:HG2	2.12	0.49
11:CL:21:PRO:C	11:CL:23:LEU:H	2.16	0.49
11:CL:41:PRO:HA	11:CL:88:ASP:O	2.12	0.49
12:CM:69:ARG:CA	12:CM:69:ARG:HH11	2.25	0.49
20:CO:48:LYS:C	20:CO:50:HIS:H	2.15	0.49
17:CT:64:GLY:C	17:CT:66:ILE:H	2.16	0.49
36:D2:45:SER:O	36:D2:46:LYS:CB	2.55	0.49
23:DB:1229:C:H2'	23:DB:1230:A:C8	2.47	0.49
23:DB:1328:A:H2'	23:DB:1330:C:C4	2.47	0.49
23:DB:2031:A:H5'	23:DB:2031:A:C8	2.46	0.49
23:DB:2066:C:O2'	23:DB:2067:G:H5'	2.11	0.49
23:DB:2344:U:H4'	23:DB:2345:G:OP1	2.12	0.49
23:DB:234:U:O2'	23:DB:235:U:H5'	2.13	0.49
23:DB:2630:G:O2'	23:DB:2631:G:H5'	2.13	0.49
23:DB:2813:A:H2'	23:DB:2814:A:H8	1.77	0.49
23:DB:413:C:H2'	23:DB:414:C:C6	2.48	0.49
23:DB:564:C:H1'	44:DQ:36:GLN:OE1	2.13	0.49
23:DB:639:U:H2'	23:DB:640:C:H6	1.77	0.49
23:DB:660:C:H2'	23:DB:661:A:H8	1.77	0.49
25:DC:123:ILE:HG13	25:DC:123:ILE:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:106:LYS:HB3	26:DD:206:ALA:CB	2.42	0.49
47:DF:167:ALA:O	47:DF:170:ALA:HB3	2.13	0.49
47:DF:30:VAL:HG21	47:DF:96:TRP:HE1	1.78	0.49
48:DG:145:ALA:O	48:DG:149:ALA:N	2.44	0.49
48:DG:167:VAL:HG23	48:DG:168:VAL:H	1.77	0.49
24:DI:128:ILE:HA	24:DI:131:THR:HG23	1.94	0.49
41:DJ:96:ARG:N	41:DJ:97:PRO:HD3	2.28	0.49
27:DK:105:ARG:N	27:DK:105:ARG:HD3	2.20	0.49
42:DN:103:ARG:HB3	42:DN:106:ASP:O	2.12	0.49
52:DW:55:ASP:C	52:DW:57:THR:H	2.16	0.49
39:DX:24:GLU:O	39:DX:28:LEU:HG	2.12	0.49
1:AA:1179:A:H2'	1:AA:1180:A:C8	2.47	0.49
1:AA:203:G:H1'	1:AA:465:A:N6	2.26	0.49
1:AA:236:A:O2'	1:AA:237:G:H5'	2.12	0.49
1:AA:284:C:H2'	1:AA:285:C:H6	1.77	0.49
1:AA:378:G:O2'	1:AA:379:C:H5'	2.13	0.49
1:AA:397:A:N3	1:AA:397:A:H3'	2.27	0.49
1:AA:462:G:H5'	1:AA:463:U:P	2.53	0.49
1:AA:82:G:N1	1:AA:88:U:H1'	2.28	0.49
4:AE:92:ARG:O	4:AE:126:ALA:HB1	2.12	0.49
4:AE:21:SER:HB3	4:AE:28:ARG:NH2	2.28	0.49
6:AG:69:ARG:HB3	6:AG:95:ARG:HD3	1.94	0.49
7:AH:77:VAL:HG23	7:AH:126:CYS:HA	1.94	0.49
10:AK:95:THR:HG23	10:AK:96:ILE:H	1.78	0.49
21:AN:5:MET:SD	21:AN:8:ARG:HD3	2.53	0.49
15:AR:41:SER:HB3	15:AR:51:GLN:CD	2.33	0.49
1:AA:673:A:H1'	15:AR:63:TYR:HE1	1.76	0.49
16:AS:10:ILE:HG21	16:AS:40:PHE:CZ	2.46	0.49
23:BB:1076:C:H4'	24:BI:94:LYS:CE	2.42	0.49
23:BB:1324:G:H1'	23:BB:1616:A:C6	2.48	0.49
23:BB:1778:U:H2'	23:BB:1784:A:N6	2.27	0.49
23:BB:1945:G:H2'	23:BB:1946:U:C6	2.48	0.49
23:BB:2201:G:H2'	23:BB:2202:U:C6	2.47	0.49
23:BB:2393:U:O2'	23:BB:2394:C:H5'	2.13	0.49
23:BB:2675:A:N1	23:BB:2732:G:O6	2.46	0.49
23:BB:267:C:H2'	23:BB:268:C:H6	1.78	0.49
23:BB:2709:G:H2'	23:BB:2710:C:C6	2.48	0.49
23:BB:38:A:H1'	29:BE:43:THR:HB	1.93	0.49
23:BB:458:G:H22	23:BB:469:G:H2'	1.78	0.49
23:BB:608:A:H2'	23:BB:609:A:H8	1.76	0.49
23:BB:768:G:H2'	23:BB:769:U:O4'	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:77:ILE:HG13	29:BE:78:TRP:HE3	1.78	0.49
47:BF:105:ILE:HA	47:BF:108:PRO:HB2	1.94	0.49
41:BJ:96:ARG:N	41:BJ:97:PRO:HD3	2.27	0.49
27:BK:61:VAL:HG11	27:BK:112:PHE:CZ	2.48	0.49
27:BK:66:LYS:HD2	27:BK:81:GLY:N	2.27	0.49
27:BK:70:ARG:CB	27:BK:76:VAL:HG22	2.42	0.49
42:BN:83:LEU:HD21	42:BN:115:LEU:HD11	1.95	0.49
44:BQ:63:ARG:HH22	44:BQ:96:ASP:CA	2.26	0.49
49:BR:24:LYS:HD3	49:BR:24:LYS:O	2.13	0.49
37:BL:23:ILE:HG13	49:BR:82:HIS:ND1	2.28	0.49
50:BT:85:VAL:C	50:BT:86:THR:HG23	2.32	0.49
50:BT:8:LEU:HD22	50:BT:46:ALA:HA	1.93	0.49
46:BU:53:GLN:O	46:BU:53:GLN:HG2	2.12	0.49
52:BW:69:GLU:HG3	52:BW:80:SER:HG	1.77	0.49
1:CA:1033:G:H2'	1:CA:1034:G:O4'	2.13	0.49
1:CA:1319:A:OP1	16:CS:4:LEU:HD11	2.13	0.49
1:CA:1505:G:H5''	1:CA:1506:U:O5'	2.12	0.49
1:CA:405:U:C5	3:CD:4:LEU:HD21	2.47	0.49
1:CA:779:C:H2'	1:CA:780:A:O4'	2.12	0.49
1:CA:792:A:H1'	1:CA:794:A:N7	2.27	0.49
1:CA:920:U:H2'	1:CA:921:U:H6	1.72	0.49
18:CB:205:ALA:O	18:CB:209:VAL:HG22	2.12	0.49
18:CB:55:GLU:HA	18:CB:58:LYS:HB3	1.93	0.49
3:CD:29:THR:HG22	3:CD:30:LYS:HD3	1.95	0.49
4:CE:46:GLY:CA	4:CE:70:MET:HA	2.42	0.49
12:CM:9:PRO:HB2	12:CM:17:ALA:HB1	1.95	0.49
22:DA:43:C:O2'	47:DF:91:ARG:HD2	2.12	0.49
23:DB:1168:G:H2'	23:DB:1169:A:H8	1.73	0.49
23:DB:1196:C:H2'	23:DB:1197:G:H8	1.78	0.49
23:DB:1569:A:H2'	23:DB:1570:A:C8	2.47	0.49
23:DB:1958:C:O2'	23:DB:1959:G:H5'	2.13	0.49
23:DB:1997:C:H2'	23:DB:1998:A:H8	1.77	0.49
23:DB:2543:G:H2'	23:DB:2544:G:H8	1.76	0.49
23:DB:2543:G:H4'	23:DB:2645:G:N2	2.28	0.49
23:DB:2551:C:H2'	23:DB:2552:U:O4'	2.13	0.49
23:DB:2722:G:H2'	23:DB:2723:C:C6	2.48	0.49
23:DB:341:C:H2'	23:DB:342:A:H8	1.77	0.49
23:DB:768:G:H2'	23:DB:769:U:O4'	2.12	0.49
23:DB:866:A:H61	23:DB:913:U:C4'	2.25	0.49
23:DB:925:A:O2'	23:DB:926:G:H5'	2.12	0.49
29:DE:134:LEU:HD23	29:DE:161:ALA:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DH:103:VAL:CG2	40:DH:108:VAL:HB	2.42	0.49
42:DN:12:ARG:HG3	42:DN:13:ASN:N	2.28	0.49
28:DP:4:ILE:O	28:DP:6:GLN:N	2.40	0.49
49:DR:24:LYS:O	49:DR:24:LYS:HD3	2.13	0.49
1:AA:1123:U:O2'	1:AA:1124:G:H5'	2.13	0.49
1:AA:1289:A:H2'	1:AA:1290:G:H5'	1.93	0.49
1:AA:1303:C:H2'	1:AA:1304:G:O4'	2.13	0.49
1:AA:1468:A:O2'	1:AA:1469:C:H5'	2.13	0.49
1:AA:426:U:H2'	1:AA:427:U:C6	2.48	0.49
1:AA:532:A:H62	2:AC:191:THR:HB	1.77	0.49
1:AA:911:U:H2'	1:AA:912:C:C6	2.48	0.49
18:AB:125:PHE:CD2	18:AB:126:ASP:N	2.81	0.49
6:AG:62:GLU:O	6:AG:66:GLU:HG3	2.13	0.49
9:AJ:34:ALA:HB1	9:AJ:76:ILE:HB	1.95	0.49
11:AL:33:CYS:HA	11:AL:54:VAL:HA	1.95	0.49
23:BB:1439:A:H5'	23:BB:1440:U:OP2	2.12	0.49
23:BB:1805:A:O2'	23:BB:1806:C:H5'	2.13	0.49
23:BB:1847:A:H4'	23:BB:1848:A:H8	1.78	0.49
23:BB:2327:A:H2'	23:BB:2328:A:C8	2.47	0.49
23:BB:433:C:O2'	23:BB:434:U:H5'	2.13	0.49
23:BB:849:A:H2'	23:BB:850:U:C6	2.48	0.49
23:BB:854:C:O2'	23:BB:855:G:H5'	2.12	0.49
25:BC:242:HIS:O	25:BC:244:VAL:HG13	2.12	0.49
23:BB:2772:C:H4'	26:BD:171:THR:CG2	2.42	0.49
26:BD:186:LEU:HD21	28:BP:3:ILE:HD11	1.95	0.49
29:BE:126:VAL:HG22	29:BE:127:GLU:N	2.22	0.49
29:BE:134:LEU:HD23	29:BE:161:ALA:HB2	1.93	0.49
48:BG:84:LYS:HG2	48:BG:85:LYS:N	2.23	0.49
40:BH:124:THR:HG22	40:BH:125:THR:H	1.78	0.49
24:BI:56:VAL:CG2	24:BI:68:PHE:HB2	2.43	0.49
24:BI:74:PRO:O	24:BI:77:VAL:HG22	2.12	0.49
27:BK:64:ARG:NH2	28:BP:67:GLU:HG3	2.27	0.49
42:BN:72:ASP:C	42:BN:74:GLU:H	2.16	0.49
23:BB:2376:A:N6	43:BO:94:ARG:HD2	2.21	0.49
28:BP:63:ILE:HA	28:BP:67:GLU:O	2.13	0.49
49:BR:29:THR:O	49:BR:63:VAL:HG23	2.12	0.49
45:BS:26:GLY:O	45:BS:28:LYS:N	2.46	0.49
39:BX:52:ARG:O	39:BX:56:LEU:HD23	2.13	0.49
30:BY:10:ARG:HB3	30:BY:10:ARG:NH2	2.27	0.49
30:BY:11:SER:OG	30:BY:13:ILE:HG13	2.12	0.49
30:BY:6:ILE:HG21	30:BY:47:ILE:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1432:G:H1'	1:CA:1468:A:H61	1.77	0.49
1:CA:781:A:O2'	1:CA:1522:U:O2	2.30	0.49
1:CA:397:A:H3'	1:CA:397:A:N3	2.27	0.49
1:CA:658:C:O2'	1:CA:659:U:H5'	2.12	0.49
1:CA:814:A:O2'	1:CA:815:A:H3'	2.13	0.49
2:CC:154:GLY:HA2	2:CC:163:ARG:O	2.13	0.49
4:CE:97:PRO:HA	4:CE:122:VAL:HG12	1.93	0.49
5:CF:97:THR:HB	5:CF:98:GLU:OE1	2.13	0.49
4:CE:156:ARG:HH12	7:CH:42:GLU:HB3	1.76	0.49
8:CI:11:ARG:HA	8:CI:105:ARG:NH2	2.28	0.49
11:CL:89:LEU:HD22	11:CL:89:LEU:N	2.28	0.49
23:DB:1119:U:H2'	23:DB:1120:G:C8	2.44	0.49
23:DB:1026:G:OP2	23:DB:1134:A:H1'	2.13	0.49
23:DB:1230:A:H2'	23:DB:1231:U:H6	1.73	0.49
23:DB:1237:A:N3	23:DB:1237:A:H2'	2.27	0.49
23:DB:1251:C:OP2	44:DQ:9:ALA:HB2	2.13	0.49
23:DB:2291:U:H2'	23:DB:2292:U:H6	1.76	0.49
23:DB:2415:G:H2'	23:DB:2416:C:C6	2.48	0.49
23:DB:2665:A:O2'	23:DB:2666:C:H5'	2.12	0.49
23:DB:41:C:O2'	23:DB:42:A:H5'	2.12	0.49
23:DB:492:A:H2'	23:DB:493:G:O4'	2.13	0.49
23:DB:526:A:N6	23:DB:2626:C:H4'	2.28	0.49
23:DB:593:U:H2'	23:DB:594:U:H6	1.74	0.49
23:DB:659:G:H4'	29:DE:95:LYS:HB3	1.95	0.49
23:DB:680:C:H2'	23:DB:681:G:C8	2.47	0.49
23:DB:863:A:H2'	23:DB:864:G:C8	2.46	0.49
25:DC:42:ARG:CZ	25:DC:48:ILE:HD11	2.43	0.49
23:DB:1655:A:H5'	26:DD:118:PHE:HB2	1.95	0.49
26:DD:154:LYS:HG2	26:DD:155:VAL:N	2.26	0.49
26:DD:58:ASN:OD1	26:DD:59:ARG:HG3	2.13	0.49
47:DF:102:LEU:O	47:DF:103:ILE:HB	2.13	0.49
47:DF:127:TYR:O	47:DF:155:ILE:HB	2.12	0.49
47:DF:169:LEU:HB3	47:DF:174:PHE:CD1	2.48	0.49
48:DG:17:LYS:HG3	48:DG:19:ASN:OD1	2.12	0.49
48:DG:72:ASN:O	48:DG:76:ILE:HG12	2.13	0.49
40:DH:135:HIS:HD2	40:DH:138:VAL:HG23	1.77	0.49
41:DJ:88:THR:HG22	41:DJ:91:GLU:OE1	2.13	0.49
42:DN:35:LYS:HA	42:DN:111:ALA:O	2.12	0.49
42:DN:96:ARG:NE	42:DN:116:VAL:HA	2.28	0.49
44:DQ:88:GLU:HG2	49:DR:49:ILE:O	2.13	0.49
52:DW:9:THR:HG23	52:DW:10:ARG:HD3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1053:G:C6	1:AA:1199:U:H2'	2.48	0.49
1:AA:1060:U:H2'	1:AA:1061:G:C8	2.47	0.49
1:AA:1132:C:H2'	1:AA:1133:G:C8	2.48	0.49
1:AA:1486:G:C2'	1:AA:1487:G:O4'	2.61	0.49
1:AA:237:G:H5''	14:AQ:26:ARG:NH2	2.28	0.49
1:AA:255:G:H1'	14:AQ:17:GLU:OE2	2.13	0.49
1:AA:420:U:H2'	1:AA:422:C:C4	2.48	0.49
1:AA:524:G:H2'	1:AA:525:C:H6	1.77	0.49
1:AA:539:A:H2'	1:AA:540:G:H8	1.78	0.49
1:AA:821:G:H2'	1:AA:822:U:H6	1.77	0.49
18:AB:52:ALA:O	18:AB:56:LEU:HD13	2.12	0.49
2:AC:96:VAL:CB	2:AC:97:PRO:HD2	2.41	0.49
4:AE:96:GLN:HB3	4:AE:123:LEU:CD1	2.42	0.49
6:AG:11:ILE:HG12	6:AG:24:LYS:HE2	1.94	0.49
9:AJ:88:MET:HA	9:AJ:91:ASP:OD1	2.13	0.49
10:AK:36:ARG:HG3	10:AK:36:ARG:HH11	1.78	0.49
12:AM:92:ARG:NH1	16:AS:79:TYR:HE2	2.10	0.49
36:B2:44:VAL:O	36:B2:45:SER:C	2.51	0.49
23:BB:136:G:O5'	23:BB:136:G:H8	1.96	0.49
23:BB:1477:A:H2'	23:BB:1478:G:O4'	2.13	0.49
23:BB:1681:G:H21	23:BB:1762:A:H3'	1.78	0.49
23:BB:2340:A:H2'	23:BB:2341:G:H8	1.78	0.49
23:BB:2543:G:H2'	23:BB:2544:G:H8	1.75	0.49
23:BB:2579:C:O5'	23:BB:2579:C:H6	1.95	0.49
23:BB:409:G:H2'	23:BB:410:G:C8	2.48	0.49
23:BB:438:G:H2'	23:BB:439:A:C8	2.47	0.49
23:BB:445:C:H2'	23:BB:446:G:O4'	2.13	0.49
23:BB:49:A:OP1	23:BB:51:G:H5'	2.13	0.49
23:BB:627:A:OP1	23:BB:627:A:H2'	2.13	0.49
26:BD:123:LYS:HB3	26:BD:165:MET:HE3	1.94	0.49
26:BD:148:GLN:CG	26:BD:152:PRO:HG2	2.43	0.49
26:BD:38:LYS:HB2	26:BD:47:ALA:HB3	1.95	0.49
47:BF:91:ARG:C	47:BF:95:MET:HB2	2.33	0.49
40:BH:18:GLN:HE21	40:BH:39:ALA:CB	2.26	0.49
42:BN:58:ASP:O	42:BN:59:SER:HB3	2.12	0.49
44:BQ:60:TRP:C	44:BQ:64:ILE:HG12	2.32	0.49
52:BW:35:ILE:HG12	52:BW:35:ILE:O	2.12	0.49
39:BX:23:ARG:HA	39:BX:27:ASN:H	1.77	0.49
1:CA:404:G:O2'	1:CA:405:U:H5'	2.12	0.49
1:CA:559:A:H4'	1:CA:560:A:H3'	1.95	0.49
1:CA:656:G:H2'	1:CA:657:U:H6	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:73:C:C2'	1:CA:74:A:H5'	2.43	0.49
18:CB:42:LEU:HA	18:CB:45:THR:OG1	2.13	0.49
18:CB:80:LYS:HG3	18:CB:81:ASP:N	2.27	0.49
2:CC:106:ARG:N	2:CC:106:ARG:NE	2.54	0.49
21:CN:19:TYR:C	21:CN:21:ALA:H	2.15	0.49
13:CP:23:ASP:CG	13:CP:25:ARG:HE	2.16	0.49
14:CQ:18:LYS:H	14:CQ:50:ASN:ND2	2.10	0.49
14:CQ:20:ILE:HG12	14:CQ:52:CYS:SG	2.52	0.49
16:CS:39:ILE:CD1	16:CS:68:HIS:H	2.25	0.49
23:DB:1526:C:H2'	23:DB:1527:G:O4'	2.12	0.49
23:DB:2094:A:H2'	23:DB:2095:A:H8	1.77	0.49
23:DB:2487:G:H2'	23:DB:2488:G:C8	2.48	0.49
23:DB:2488:G:H2'	23:DB:2489:U:C6	2.48	0.49
23:DB:2686:G:H2'	23:DB:2687:U:C6	2.48	0.49
23:DB:408:G:O2'	23:DB:409:G:H5'	2.13	0.49
23:DB:693:A:H2'	23:DB:694:U:H6	1.78	0.49
23:DB:917:A:H5''	23:DB:2268:A:N6	2.27	0.49
23:DB:992:C:H2'	23:DB:993:G:H8	1.77	0.49
29:DE:145:ASP:H	29:DE:166:LYS:CB	2.26	0.49
29:DE:46:GLN:HG3	29:DE:87:ALA:HB3	1.95	0.49
48:DG:59:ASP:O	48:DG:63:GLN:HB2	2.13	0.49
40:DH:135:HIS:CD2	40:DH:138:VAL:HG23	2.48	0.49
37:DL:109:LYS:HB2	37:DL:111:ILE:HD11	1.94	0.49
43:DO:52:SER:C	43:DO:54:VAL:H	2.16	0.49
49:DR:31:GLU:H	49:DR:63:VAL:CG2	2.25	0.49
45:DS:55:ILE:HD12	45:DS:107:VAL:HG21	1.94	0.49
23:DB:751:A:H5'	45:DS:90:LYS:HA	1.95	0.49
52:DW:43:LYS:HE3	52:DW:68:PHE:HE1	1.78	0.49
39:DX:49:ASP:O	39:DX:53:VAL:HG23	2.13	0.49
1:AA:370:C:H2'	1:AA:371:A:C8	2.47	0.48
1:AA:373:A:P	1:AA:373:A:H3'	2.53	0.48
18:AB:134:LEU:HA	18:AB:137:THR:OG1	2.12	0.48
18:AB:136:ARG:O	18:AB:139:GLU:HB3	2.13	0.48
2:AC:155:ARG:NH1	2:AC:160:GLU:HA	2.28	0.48
1:AA:406:G:N2	3:AD:115:GLN:HE22	2.10	0.48
3:AD:117:VAL:O	3:AD:130:ASN:HA	2.12	0.48
4:AE:39:GLY:HA3	4:AE:116:VAL:O	2.13	0.48
7:AH:10:LEU:HA	7:AH:74:ILE:HD11	1.95	0.48
17:AT:54:GLN:N	17:AT:55:PRO:HD2	2.27	0.48
31:B0:32:THR:OG1	31:B0:50:GLY:HA2	2.12	0.48
32:B4:15:LYS:O	32:B4:16:ILE:HB	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:BA:35:C:H2'	22:BA:36:C:O4'	2.13	0.48
23:BB:1306:C:O2'	23:BB:1307:A:H5'	2.13	0.48
23:BB:1544:A:H2'	23:BB:1545:A:C8	2.48	0.48
23:BB:1593:A:H2'	23:BB:1594:U:H6	1.78	0.48
23:BB:1711:A:H2'	23:BB:1712:U:C6	2.48	0.48
23:BB:1993:U:H4'	26:BD:133:THR:HG21	1.95	0.48
23:BB:688:U:H2'	23:BB:689:A:H8	1.78	0.48
26:BD:118:PHE:O	26:BD:119:ALA:CB	2.61	0.48
26:BD:1:MET:SD	26:BD:2:ILE:N	2.85	0.48
26:BD:106:LYS:HB3	26:BD:206:ALA:CB	2.43	0.48
29:BE:131:THR:HG22	29:BE:161:ALA:H	1.78	0.48
47:BF:141:ASP:CB	47:BF:144:LYS:HB2	2.43	0.48
48:BG:145:ALA:O	48:BG:149:ALA:N	2.44	0.48
24:BI:21:PRO:CB	24:BI:22:PRO:HD3	2.39	0.48
41:BJ:20:ALA:HB2	41:BJ:28:LEU:HD13	1.94	0.48
27:BK:17:ARG:HB3	27:BK:45:GLU:HG3	1.93	0.48
44:BQ:75:TYR:O	44:BQ:79:ILE:HG22	2.13	0.48
45:BS:45:VAL:HG23	45:BS:46:LEU:N	2.28	0.48
46:BU:73:ASN:OD1	46:BU:75:ALA:HB3	2.13	0.48
1:CA:1237:C:H2'	1:CA:1336:C:C5	2.48	0.48
1:CA:160:A:H2'	1:CA:161:A:O4'	2.14	0.48
1:CA:755:G:O2'	1:CA:756:C:H5'	2.13	0.48
1:CA:922:G:H2'	1:CA:923:A:H8	1.78	0.48
3:CD:117:VAL:O	3:CD:130:ASN:HA	2.13	0.48
4:CE:37:VAL:HG12	4:CE:47:PHE:HB2	1.94	0.48
7:CH:124:ILE:C	7:CH:125:ILE:HD12	2.34	0.48
7:CH:37:ASN:O	7:CH:41:GLU:HG2	2.13	0.48
8:CI:52:GLU:O	8:CI:53:LEU:HD22	2.12	0.48
9:CJ:7:ARG:HG2	9:CJ:75:ASP:OD2	2.13	0.48
9:CJ:9:ARG:HB2	9:CJ:99:GLN:CB	2.43	0.48
10:CK:65:ALA:O	10:CK:68:ARG:HB3	2.12	0.48
12:CM:13:HIS:HB3	12:CM:40:GLU:O	2.12	0.48
33:D1:36:LYS:HG3	33:D1:46:VAL:O	2.13	0.48
33:D1:49:LYS:HG2	33:D1:50:GLU:N	2.16	0.48
22:DA:96:G:O2'	22:DA:97:C:H5'	2.13	0.48
23:DB:1098:A:H2'	24:DI:4:VAL:C	2.34	0.48
23:DB:1550:C:H2'	23:DB:1551:A:C8	2.48	0.48
23:DB:15:G:O2'	23:DB:16:C:H5'	2.13	0.48
23:DB:1947:C:O2'	23:DB:1948:G:H5'	2.12	0.48
23:DB:2519:U:C6	23:DB:2542:A:N6	2.81	0.48
23:DB:2553:G:H2'	23:DB:2554:U:O4'	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2633:G:H2'	23:DB:2634:A:O4'	2.13	0.48
23:DB:2636:C:O2'	23:DB:2637:U:H5'	2.13	0.48
23:DB:2654:A:N1	23:DB:2665:A:H5''	2.28	0.48
23:DB:2897:U:H2'	23:DB:2898:U:C6	2.48	0.48
23:DB:560:C:H3'	23:DB:561:G:H8	1.78	0.48
23:DB:714:U:O2	23:DB:716:A:H3'	2.13	0.48
23:DB:796:C:H2'	23:DB:797:G:C8	2.48	0.48
23:DB:79:C:O2'	23:DB:346:A:H1'	2.13	0.48
25:DC:155:ARG:HB3	25:DC:155:ARG:HH11	1.77	0.48
25:DC:209:ALA:HA	25:DC:212:TRP:NE1	2.28	0.48
25:DC:242:HIS:O	25:DC:244:VAL:HG13	2.13	0.48
25:DC:250:GLN:O	25:DC:250:GLN:HG2	2.13	0.48
29:DE:62:GLN:HE21	29:DE:62:GLN:HB2	1.52	0.48
23:DB:2312:U:H4'	47:DF:84:ILE:HG21	1.94	0.48
40:DH:70:GLU:OE1	40:DH:71:LYS:HD2	2.12	0.48
24:DI:79:LEU:HD12	24:DI:135:MET:SD	2.52	0.48
41:DJ:15:TRP:HB3	41:DJ:137:PRO:HB3	1.94	0.48
27:DK:2:ILE:HG22	27:DK:32:TYR:HB3	1.94	0.48
37:DL:116:VAL:HG22	37:DL:117:THR:N	2.28	0.48
37:DL:134:ALA:O	37:DL:137:ALA:HB3	2.13	0.48
42:DN:72:ASP:OD2	42:DN:74:GLU:HB3	2.13	0.48
43:DO:105:ALA:O	43:DO:106:LEU:HG	2.13	0.48
43:DO:47:VAL:HG12	43:DO:48:LEU:N	2.22	0.48
49:DR:39:LEU:HD23	49:DR:39:LEU:N	2.28	0.48
45:DS:28:LYS:HB2	45:DS:31:GLN:HB2	1.95	0.48
50:DT:50:LEU:C	50:DT:52:GLU:H	2.16	0.48
46:DU:53:GLN:HG2	46:DU:53:GLN:O	2.13	0.48
35:DV:14:LYS:CE	35:DV:18:ARG:HH21	2.24	0.48
35:DV:65:VAL:O	35:DV:67:GLY:N	2.44	0.48
51:DZ:5:CYS:HB3	51:DZ:9:GLY:N	2.27	0.48
1:AA:1039:G:H2'	1:AA:1040:U:C6	2.48	0.48
1:AA:1040:U:H2'	1:AA:1041:G:C8	2.47	0.48
1:AA:1067:A:H4'	1:AA:1068:G:O5'	2.13	0.48
1:AA:1273:C:H2'	1:AA:1274:A:O4'	2.13	0.48
1:AA:1355:G:H2'	1:AA:1356:G:C8	2.48	0.48
1:AA:363:A:P	11:AL:57:THR:HG21	2.54	0.48
1:AA:845:A:H3'	1:AA:846:G:H8	1.78	0.48
1:AA:985:C:H2'	1:AA:986:U:C6	2.49	0.48
1:AA:986:U:C1'	16:AS:54:ARG:HB3	2.43	0.48
11:AL:30:ARG:O	11:AL:57:THR:HG23	2.13	0.48
11:AL:89:LEU:HD22	11:AL:89:LEU:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:BA:83:G:OP1	30:BY:16:LEU:HD23	2.14	0.48
23:BB:1312:U:H4'	23:BB:1313:U:O5'	2.13	0.48
23:BB:1623:G:O2'	23:BB:1624:U:H5'	2.13	0.48
23:BB:1747:U:H2'	23:BB:1748:C:H6	1.77	0.48
23:BB:1808:A:O2'	23:BB:1809:A:H5'	2.13	0.48
23:BB:2243:U:H2'	23:BB:2244:U:C6	2.48	0.48
23:BB:2389:G:C5'	23:BB:2390:U:H5'	2.42	0.48
23:BB:2419:U:OP2	34:B3:32:LEU:HD13	2.13	0.48
23:BB:2551:C:H2'	23:BB:2552:U:O4'	2.14	0.48
23:BB:2853:C:H2'	23:BB:2854:G:H8	1.77	0.48
42:BN:49:GLU:HA	42:BN:94:TYR:HD2	1.78	0.48
42:BN:72:ASP:OD2	42:BN:74:GLU:HB3	2.14	0.48
28:BP:16:VAL:HG23	28:BP:16:VAL:O	2.13	0.48
49:BR:49:ILE:HG12	49:BR:53:PHE:HA	1.95	0.48
46:BU:48:VAL:O	46:BU:50:ALA:N	2.46	0.48
46:BU:3:LYS:CB	46:BU:82:VAL:HG21	2.38	0.48
23:BB:2355:G:H4'	52:BW:20:LEU:CD1	2.42	0.48
1:CA:1328:C:H5''	12:CM:27:THR:CG2	2.42	0.48
1:CA:375:U:OP1	13:CP:70:ARG:HB2	2.13	0.48
1:CA:651:C:H2'	1:CA:652:U:C6	2.48	0.48
1:CA:853:C:O2'	1:CA:854:U:H5'	2.13	0.48
4:CE:81:GLN:HG2	4:CE:148:SER:HA	1.95	0.48
8:CI:109:GLN:CD	8:CI:110:VAL:H	2.16	0.48
8:CI:48:ARG:O	8:CI:52:GLU:N	2.44	0.48
9:CJ:52:LEU:CD1	9:CJ:52:LEU:H	2.22	0.48
21:CN:10:VAL:HB	21:CN:11:LYS:NZ	2.28	0.48
16:CS:33:TRP:C	16:CS:35:ARG:H	2.17	0.48
23:DB:1099:G:O2'	23:DB:1100:C:H5'	2.13	0.48
23:DB:1259:G:H2'	23:DB:1260:A:H8	1.78	0.48
23:DB:1677:A:H2'	23:DB:1678:A:C8	2.48	0.48
23:DB:17:G:H2'	23:DB:18:U:H6	1.79	0.48
23:DB:195:A:H1'	23:DB:250:G:N2	2.28	0.48
23:DB:2428:G:H5''	23:DB:2429:G:OP1	2.14	0.48
23:DB:2645:G:H5''	23:DB:2732:G:C8	2.44	0.48
23:DB:721:A:H2'	23:DB:722:A:C8	2.47	0.48
23:DB:942:G:H2'	23:DB:943:A:H8	1.78	0.48
26:DD:106:LYS:HB3	26:DD:206:ALA:HB3	1.95	0.48
23:DB:2680:U:OP2	26:DD:114:LYS:HB3	2.12	0.48
29:DE:88:ARG:O	29:DE:90:GLN:HG3	2.13	0.48
47:DF:90:LEU:C	47:DF:91:ARG:HD3	2.33	0.48
48:DG:106:LEU:HD13	48:DG:151:ARG:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DK:19:VAL:HB	27:DK:41:ILE:CD1	2.43	0.48
23:DB:1952:A:H5'	27:DK:42:THR:HG21	1.95	0.48
27:DK:92:GLU:HB2	27:DK:93:GLN:HE21	1.77	0.48
42:DN:83:LEU:HD21	42:DN:115:LEU:HD11	1.95	0.48
43:DO:26:LEU:HG	43:DO:26:LEU:O	2.12	0.48
28:DP:47:ILE:HA	28:DP:96:LEU:HB2	1.94	0.48
44:DQ:59:LEU:HD13	44:DQ:60:TRP:N	2.28	0.48
51:DZ:14:THR:HA	51:DZ:28:ARG:HA	1.95	0.48
1:AA:1102:A:OP2	1:AA:1102:A:H8	1.95	0.48
1:AA:1231:G:H2'	1:AA:1232:U:H6	1.75	0.48
1:AA:1499:A:O2'	1:AA:1500:A:H5'	2.12	0.48
1:AA:254:G:H4'	14:AQ:19:SER:OG	2.13	0.48
1:AA:502:A:H4'	1:AA:550:G:H4'	1.95	0.48
1:AA:613:C:H2'	1:AA:614:C:H6	1.78	0.48
1:AA:724:G:H2'	1:AA:725:G:C8	2.45	0.48
6:AG:150:PHE:H	10:AK:55:ARG:NH2	2.11	0.48
21:AN:20:PHE:O	21:AN:21:ALA:C	2.51	0.48
21:AN:26:LEU:HD23	21:AN:27:LYS:H	1.78	0.48
21:AN:45:LEU:HD23	21:AN:45:LEU:O	2.14	0.48
21:AN:86:ALA:HB3	21:AN:92:ILE:HD11	1.94	0.48
14:AQ:46:HIS:HB2	14:AQ:66:LEU:HD13	1.95	0.48
23:BB:1059:G:H2'	23:BB:1060:U:C5	2.48	0.48
23:BB:1361:G:H2'	23:BB:1362:C:C6	2.49	0.48
23:BB:150:U:O2'	23:BB:151:C:H5'	2.12	0.48
23:BB:172:A:H2'	23:BB:173:A:C8	2.48	0.48
23:BB:1859:U:O5'	23:BB:1859:U:H6	1.97	0.48
23:BB:2530:A:H2'	23:BB:2531:A:H5''	1.95	0.48
23:BB:2656:U:H2'	23:BB:2657:A:H8	1.77	0.48
23:BB:2783:U:H2'	23:BB:2784:U:C6	2.48	0.48
23:BB:416:U:H2'	23:BB:417:C:C6	2.48	0.48
25:BC:9:SER:C	25:BC:11:GLY:H	2.16	0.48
26:BD:11:MET:H	26:BD:25:THR:HA	1.78	0.48
29:BE:28:VAL:HG23	29:BE:29:HIS:N	2.28	0.48
47:BF:111:ARG:HE	47:BF:135:ILE:CG2	2.26	0.48
48:BG:143:VAL:O	48:BG:147:LEU:HG	2.13	0.48
48:BG:33:THR:HA	48:BG:34:ARG:NH1	2.28	0.48
40:BH:90:LEU:HD23	40:BH:92:GLY:H	1.79	0.48
37:BL:76:GLU:HG3	37:BL:111:ILE:HG13	1.94	0.48
42:BN:28:LEU:N	42:BN:34:ILE:HD11	2.28	0.48
42:BN:55:ALA:HA	42:BN:80:PHE:CD1	2.48	0.48
49:BR:2:TYR:N	49:BR:42:ALA:HB2	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BS:28:LYS:HB2	45:BS:31:GLN:HB2	1.95	0.48
46:BU:35:VAL:HG23	46:BU:38:ILE:HG21	1.95	0.48
30:BY:53:MET:HA	30:BY:53:MET:HE2	1.95	0.48
1:CA:1053:G:H4'	1:CA:1054:C:H5'	1.95	0.48
1:CA:191:G:H8	1:CA:191:G:OP2	1.96	0.48
1:CA:65:A:C2	1:CA:381:C:H2'	2.49	0.48
1:CA:706:A:H4'	10:CK:30:ILE:HD11	1.95	0.48
1:CA:83:C:H1'	1:CA:84:U:C5	2.48	0.48
1:CA:22:G:H4'	1:CA:885:G:C8	2.49	0.48
1:CA:895:G:H2'	1:CA:896:C:C6	2.48	0.48
1:CA:939:G:H4'	6:CG:101:ARG:HH22	1.78	0.48
18:CB:116:LEU:HD23	18:CB:119:GLN:HG3	1.95	0.48
18:CB:57:ASN:OD1	18:CB:57:ASN:N	2.47	0.48
18:CB:95:TRP:CD1	18:CB:171:ALA:HB2	2.47	0.48
2:CC:34:SER:OG	2:CC:58:ARG:NH1	2.46	0.48
7:CH:54:THR:HG23	7:CH:55:LYS:N	2.28	0.48
9:CJ:32:THR:OG1	9:CJ:82:LYS:HB2	2.14	0.48
33:D1:16:THR:HG21	33:D1:39:ASP:CG	2.33	0.48
23:DB:1118:C:H2'	23:DB:1119:U:O4'	2.12	0.48
23:DB:1212:G:H1'	23:DB:1236:G:N2	2.28	0.48
23:DB:1564:C:H2'	23:DB:1565:C:C6	2.48	0.48
23:DB:1842:G:H2'	23:DB:1843:C:C6	2.48	0.48
23:DB:2281:A:H4'	23:DB:2389:G:N2	2.28	0.48
23:DB:2784:U:O2'	23:DB:2785:C:H5'	2.14	0.48
23:DB:2809:A:H2'	23:DB:2810:A:C8	2.48	0.48
23:DB:443:A:C5	29:DE:40:ARG:HD3	2.48	0.48
23:DB:467:G:O2'	23:DB:468:G:H5'	2.13	0.48
23:DB:955:U:P	38:DM:86:LYS:HZ1	2.35	0.48
25:DC:70:LYS:HD3	25:DC:95:TYR:CE1	2.48	0.48
26:DD:148:GLN:CG	26:DD:152:PRO:HG2	2.42	0.48
26:DD:62:LYS:HD2	26:DD:62:LYS:N	2.28	0.48
47:DF:76:PHE:HD2	47:DF:78:ILE:HD13	1.78	0.48
24:DI:19:PRO:HB2	24:DI:22:PRO:HD2	1.95	0.48
23:DB:1099:G:O5'	24:DI:4:VAL:HG12	2.13	0.48
42:DN:2:ARG:HB3	42:DN:2:ARG:NH1	2.27	0.48
44:DQ:33:VAL:HG23	44:DQ:34:ALA:N	2.28	0.48
44:DQ:94:LEU:C	44:DQ:96:ASP:H	2.15	0.48
46:DU:64:ILE:HD11	46:DU:68:ASN:HD22	1.77	0.48
1:AA:1011:C:O2'	1:AA:1012:A:H5'	2.13	0.48
1:AA:987:G:N2	1:AA:1015:G:N2	2.62	0.48
1:AA:1080:A:H5''	4:AE:20:VAL:HG11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1228:C:H5'	1:AA:1229:A:OP2	2.14	0.48
1:AA:1351:U:O2'	1:AA:1352:C:H5'	2.13	0.48
1:AA:811:C:O2'	1:AA:901:A:N1	2.46	0.48
18:AB:46:VAL:N	18:AB:47:PRO:CD	2.76	0.48
2:AC:19:SER:HB3	2:AC:21:TRP:NE1	2.28	0.48
2:AC:39:ARG:CD	2:AC:56:ILE:HD11	2.42	0.48
4:AE:46:GLY:CA	4:AE:70:MET:HA	2.40	0.48
5:AF:51:ILE:O	5:AF:51:ILE:HG23	2.14	0.48
6:AG:71:THR:O	6:AG:72:VAL:HG13	2.12	0.48
8:AI:66:VAL:O	8:AI:66:VAL:HG13	2.14	0.48
23:BB:1501:G:O2'	23:BB:1502:A:H5'	2.12	0.48
23:BB:1535:A:O2'	23:BB:1536:C:H5'	2.14	0.48
23:BB:1589:U:H2'	23:BB:1590:A:C8	2.48	0.48
23:BB:1824:G:O2'	23:BB:1825:U:H5'	2.14	0.48
23:BB:2075:U:H2'	23:BB:2238:G:N2	2.29	0.48
23:BB:2276:G:O2'	23:BB:2277:G:H5'	2.12	0.48
23:BB:2531:A:N3	23:BB:2531:A:H2'	2.29	0.48
23:BB:233:A:N6	23:BB:428:A:H61	2.11	0.48
23:BB:510:C:H2'	23:BB:511:U:O4'	2.14	0.48
23:BB:795:C:H2'	23:BB:796:C:C6	2.49	0.48
23:BB:876:C:H2'	23:BB:877:A:OP1	2.14	0.48
25:BC:70:LYS:HG3	25:BC:101:ARG:NH1	2.29	0.48
26:BD:111:GLY:N	26:BD:194:PRO:HG2	2.26	0.48
23:BB:2787:C:H4'	26:BD:62:LYS:HB3	1.96	0.48
47:BF:2:LYS:N	47:BF:2:LYS:HE3	2.28	0.48
40:BH:26:ALA:O	40:BH:27:ARG:C	2.52	0.48
41:BJ:25:LEU:O	41:BJ:27:ARG:N	2.45	0.48
38:BM:102:LEU:HD22	38:BM:102:LEU:N	2.28	0.48
38:BM:83:GLY:O	38:BM:84:LYS:HG2	2.13	0.48
42:BN:29:VAL:HG13	42:BN:83:LEU:HD22	1.95	0.48
49:BR:39:LEU:HD23	49:BR:39:LEU:N	2.29	0.48
50:BT:43:ILE:CG2	50:BT:58:VAL:HG21	2.43	0.48
23:BB:483:A:H5''	46:BU:46:LYS:HG3	1.96	0.48
23:BB:2353:G:N3	52:BW:30:VAL:HG13	2.29	0.48
51:BZ:77:LYS:O	51:BZ:78:TYR:HB3	2.13	0.48
1:CA:472:U:H2'	1:CA:473:U:C6	2.48	0.48
1:CA:499:A:H1'	1:CA:500:G:C8	2.48	0.48
1:CA:842:U:H6	1:CA:842:U:OP1	1.96	0.48
1:CA:919:A:O2'	1:CA:920:U:H5'	2.12	0.48
18:CB:118:THR:O	18:CB:122:ASP:N	2.45	0.48
18:CB:16:GLY:HA2	18:CB:40:ILE:CG1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:180:ASP:C	2:CC:181:ILE:HD12	2.34	0.48
2:CC:39:ARG:CG	2:CC:56:ILE:HD11	2.43	0.48
3:CD:167:PRO:HG2	3:CD:170:LEU:HD11	1.94	0.48
5:CF:62:MET:HG3	5:CF:64:VAL:CG2	2.42	0.48
7:CH:43:GLY:HA2	7:CH:63:LYS:NZ	2.29	0.48
12:CM:15:VAL:HG23	12:CM:16:ILE:N	2.28	0.48
12:CM:2:ARG:HB3	12:CM:6:ILE:N	2.28	0.48
12:CM:14:ALA:HB1	12:CM:33:LEU:CD2	2.43	0.48
12:CM:74:MET:HA	12:CM:77:LYS:HB3	1.95	0.48
12:CM:76:ILE:O	12:CM:79:LEU:HG	2.13	0.48
19:CU:35:GLU:HB2	19:CU:37:TYR:CZ	2.49	0.48
23:DB:1098:A:H3'	24:DI:3:LYS:CB	2.42	0.48
23:DB:125:A:H3'	23:DB:126:A:C5'	2.44	0.48
23:DB:1504:A:H2'	23:DB:1505:A:C8	2.48	0.48
23:DB:1645:G:H5''	23:DB:1646:C:H5'	1.95	0.48
23:DB:1778:U:H2'	23:DB:1784:A:N6	2.26	0.48
23:DB:1997:C:H2'	23:DB:1998:A:C8	2.48	0.48
23:DB:2010:G:H2'	23:DB:2011:U:C6	2.49	0.48
23:DB:2029:G:H2'	23:DB:2031:A:OP1	2.13	0.48
23:DB:2339:C:H2'	23:DB:2340:A:C8	2.48	0.48
23:DB:2710:C:H2'	23:DB:2711:A:H8	1.78	0.48
23:DB:271:G:N2	23:DB:367:G:H1'	2.28	0.48
23:DB:341:C:H2'	23:DB:342:A:C8	2.49	0.48
23:DB:903:C:H2'	23:DB:904:G:H8	1.79	0.48
23:DB:945:A:H4'	23:DB:945:A:OP2	2.13	0.48
26:DD:89:GLU:HG3	26:DD:94:GLN:OE1	2.13	0.48
47:DF:102:LEU:HD13	47:DF:103:ILE:N	2.28	0.48
47:DF:107:VAL:HB	47:DF:108:PRO:HD3	1.96	0.48
47:DF:111:ARG:HE	47:DF:135:ILE:CG2	2.24	0.48
48:DG:76:ILE:HB	48:DG:82:PHE:CZ	2.48	0.48
40:DH:94:ILE:O	40:DH:122:LEU:HB2	2.13	0.48
44:DQ:60:TRP:C	44:DQ:64:ILE:HG12	2.33	0.48
46:DU:62:ALA:O	46:DU:63:ALA:HB3	2.13	0.48
46:DU:82:VAL:N	46:DU:96:LYS:HZ2	2.08	0.48
35:DV:60:VAL:HG12	35:DV:61:LEU:H	1.78	0.48
52:DW:23:LYS:NZ	52:DW:24:ARG:HG3	2.28	0.48
30:DY:37:ARG:HG3	30:DY:38:GLU:OE1	2.13	0.48
1:AA:186:C:H2'	1:AA:187:G:O4'	2.13	0.48
1:AA:240:G:OP1	1:AA:240:G:H4'	2.14	0.48
1:AA:51:A:H5''	1:AA:52:C:H5''	1.96	0.48
1:AA:565:U:H3'	1:AA:566:G:H2'	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:682:G:O2'	1:AA:683:G:H5'	2.13	0.48
1:AA:696:A:O2'	1:AA:697:U:H5'	2.13	0.48
1:AA:779:C:H2'	1:AA:780:A:O4'	2.13	0.48
3:AD:18:LEU:HD23	3:AD:20:LEU:HD11	1.96	0.48
8:AI:11:ARG:H	8:AI:14:SER:CB	2.25	0.48
11:AL:106:VAL:HG23	11:AL:116:TYR:HB3	1.95	0.48
12:AM:2:ARG:HA	12:AM:8:ILE:HG23	1.93	0.48
20:AO:26:GLU:HG3	20:AO:77:ARG:HH12	1.78	0.48
14:AQ:17:GLU:O	14:AQ:18:LYS:HB2	2.14	0.48
23:BB:1039:A:H4'	35:BV:46:LYS:NZ	2.28	0.48
23:BB:138:U:H2'	23:BB:140:C:O4'	2.13	0.48
23:BB:1518:C:H2'	23:BB:1519:G:C8	2.48	0.48
23:BB:1561:C:H2'	23:BB:1562:U:C6	2.49	0.48
23:BB:2010:G:H2'	23:BB:2011:U:C6	2.48	0.48
23:BB:2305:U:O2'	47:BF:132:ARG:HA	2.12	0.48
23:BB:2417:C:O2'	23:BB:2418:A:H5'	2.13	0.48
23:BB:2676:C:H2'	23:BB:2677:G:C8	2.48	0.48
23:BB:272:A:H2'	23:BB:273:G:C8	2.47	0.48
23:BB:2849:U:H4'	23:BB:2850:A:C5'	2.43	0.48
23:BB:2869:G:H2'	23:BB:2870:C:H6	1.78	0.48
23:BB:292:U:H2'	23:BB:293:U:O4'	2.14	0.48
23:BB:350:G:H2'	23:BB:351:C:C6	2.48	0.48
23:BB:492:A:H2'	23:BB:493:G:O4'	2.13	0.48
23:BB:757:G:H2'	23:BB:758:C:H5'	1.95	0.48
25:BC:92:LEU:HD11	25:BC:100:ARG:HB2	1.95	0.48
26:BD:63:PRO:C	26:BD:65:ALA:H	2.17	0.48
47:BF:71:LYS:O	47:BF:72:SER:HB3	2.13	0.48
48:BG:1:SER:H1	48:BG:61:TRP:HE3	1.60	0.48
41:BJ:56:VAL:C	41:BJ:57:LEU:HD12	2.34	0.48
42:BN:3:HIS:O	42:BN:4:ARG:HB2	2.13	0.48
22:BA:113:C:H1'	43:BO:46:GLU:HA	1.96	0.48
28:BP:19:PHE:HE2	28:BP:83:ILE:HD11	1.77	0.48
45:BS:13:SER:OG	45:BS:14:ALA:N	2.47	0.48
52:BW:9:THR:HG23	52:BW:10:ARG:HD3	1.95	0.48
39:BX:7:ARG:HE	39:BX:7:ARG:HA	1.78	0.48
30:BY:8:GLN:HB3	30:BY:31:ILE:O	2.13	0.48
1:CA:1237:C:O2'	1:CA:1300:G:N2	2.44	0.48
1:CA:1238:A:H2	1:CA:1241:G:N3	2.12	0.48
1:CA:300:A:H2'	1:CA:301:G:O4'	2.13	0.48
1:CA:321:A:O2'	1:CA:322:C:H5'	2.12	0.48
1:CA:714:G:H2'	1:CA:715:A:C8	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:792:A:H4'	1:CA:793:U:O5'	2.13	0.48
1:CA:942:G:H21	8:CI:125:GLN:NE2	2.03	0.48
1:CA:991:U:H5	1:CA:1212:U:HO2'	1.60	0.48
6:CG:21:LEU:HD12	6:CG:100:MET:HE2	1.95	0.48
10:CK:23:HIS:HB3	10:CK:30:ILE:CG1	2.42	0.48
11:CL:80:LEU:HB3	11:CL:97:VAL:CG2	2.44	0.48
12:CM:48:SER:O	12:CM:52:ILE:HG22	2.13	0.48
21:CN:60:ARG:O	21:CN:62:ARG:HG3	2.13	0.48
13:CP:41:PRO:O	13:CP:42:ILE:HD13	2.13	0.48
13:CP:6:LEU:HB3	13:CP:17:TYR:HB3	1.94	0.48
16:CS:10:ILE:HG22	16:CS:14:LEU:CD2	2.41	0.48
1:CA:958:A:N1	16:CS:53:GLY:HA3	2.27	0.48
33:D1:8:ILE:CD1	33:D1:9:LYS:H	2.26	0.48
23:DB:1446:C:H2'	23:DB:1447:C:C6	2.49	0.48
23:DB:1464:G:H2'	23:DB:1465:G:H8	1.77	0.48
23:DB:2623:G:O2'	23:DB:2624:G:H5'	2.13	0.48
25:DC:76:VAL:HA	25:DC:114:GLN:HA	1.96	0.48
26:DD:123:LYS:HB3	26:DD:165:MET:HE3	1.95	0.48
23:DB:2578:G:N7	26:DD:145:SER:HB2	2.28	0.48
26:DD:154:LYS:H	26:DD:154:LYS:HD3	1.79	0.48
26:DD:187:LEU:HD12	26:DD:188:LEU:H	1.78	0.48
26:DD:106:LYS:CB	26:DD:206:ALA:H	2.23	0.48
29:DE:131:THR:HG22	29:DE:161:ALA:H	1.79	0.48
29:DE:136:GLN:HB2	29:DE:139:LYS:HE3	1.94	0.48
41:DJ:20:ALA:HB2	41:DJ:28:LEU:CD1	2.43	0.48
23:DB:663:G:OP1	37:DL:17:LYS:HG2	2.13	0.48
28:DP:19:PHE:CZ	28:DP:25:VAL:HG11	2.47	0.48
44:DQ:38:VAL:O	44:DQ:39:ILE:C	2.52	0.48
50:DT:21:SER:O	50:DT:25:GLU:HB2	2.13	0.48
35:DV:42:LEU:N	35:DV:42:LEU:HD23	2.29	0.48
35:DV:63:ILE:HB	35:DV:70:ILE:CG1	2.43	0.48
51:DZ:59:ILE:CD1	51:DZ:67:VAL:HG21	2.44	0.48
1:AA:1217:C:H2'	1:AA:1218:C:C6	2.48	0.48
1:AA:1404:C:H2'	1:AA:1405:G:C8	2.48	0.48
1:AA:315:A:H5''	1:AA:317:U:OP2	2.14	0.48
1:AA:585:G:O2'	1:AA:586:C:H5'	2.13	0.48
1:AA:757:U:O2'	1:AA:879:C:H1'	2.13	0.48
1:AA:756:C:H2'	1:AA:757:U:O4'	2.14	0.48
1:AA:812:G:O2'	1:AA:813:U:C6	2.65	0.48
1:AA:895:G:H2'	1:AA:896:C:H6	1.79	0.48
18:AB:18:GLN:O	18:AB:37:VAL:HG23	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:152:VAL:HG23	2:AC:165:GLU:HB3	1.96	0.48
2:AC:40:GLN:HG3	2:AC:41:TYR:H	1.77	0.48
3:AD:146:GLU:N	3:AD:146:GLU:CD	2.66	0.48
3:AD:197:HIS:ND1	3:AD:198:LEU:N	2.61	0.48
4:AE:56:PRO:HA	4:AE:59:ILE:CD1	2.44	0.48
5:AF:22:ILE:O	5:AF:26:THR:HG23	2.13	0.48
6:AG:53:SER:C	6:AG:55:LYS:H	2.16	0.48
7:AH:46:GLU:N	7:AH:63:LYS:HE3	2.29	0.48
7:AH:68:LYS:HD2	7:AH:69:ALA:H	1.79	0.48
9:AJ:44:THR:OG1	9:AJ:70:HIS:HA	2.14	0.48
10:AK:65:ALA:O	10:AK:68:ARG:HB3	2.13	0.48
17:AT:72:ALA:HA	17:AT:75:LYS:HD3	1.95	0.48
34:B3:57:VAL:C	34:B3:59:ALA:H	2.16	0.48
22:BA:116:G:H4'	43:BO:54:VAL:HG22	1.96	0.48
23:BB:1005:C:H2'	23:BB:1006:C:C6	2.48	0.48
23:BB:1222:U:P	49:BR:90:ARG:HH12	2.37	0.48
23:BB:170:U:H2'	23:BB:171:U:H6	1.77	0.48
23:BB:2246:G:H2'	23:BB:2247:A:C8	2.49	0.48
23:BB:2645:G:H5''	23:BB:2732:G:C8	2.45	0.48
23:BB:2789:C:H2'	23:BB:2790:U:H5'	1.96	0.48
23:BB:2881:U:O2'	23:BB:2882:A:H5'	2.14	0.48
23:BB:996:A:O3'	44:BQ:91:ARG:HG2	2.14	0.48
25:BC:250:GLN:O	25:BC:250:GLN:HG2	2.13	0.48
26:BD:107:VAL:N	26:BD:206:ALA:H	2.11	0.48
29:BE:126:VAL:HG13	29:BE:156:ASN:HD22	1.78	0.48
24:BI:89:SER:HB2	24:BI:136:GLY:HA3	1.94	0.48
43:BO:76:LYS:HG3	43:BO:113:ALA:HB2	1.95	0.48
43:BO:26:LEU:O	43:BO:26:LEU:HG	2.14	0.48
44:BQ:111:LYS:HE2	49:BR:50:GLY:CA	2.44	0.48
50:BT:29:THR:CG2	50:BT:86:THR:HG22	2.43	0.48
23:BB:397:U:H5''	51:BZ:32:ASN:ND2	2.28	0.48
1:CA:1095:U:H2'	1:CA:1096:C:C6	2.48	0.48
1:CA:1246:A:H2'	1:CA:1247:U:O4'	2.14	0.48
1:CA:1275:A:H2'	1:CA:1276:G:O4'	2.14	0.48
1:CA:186:C:H2'	1:CA:187:G:O4'	2.14	0.48
1:CA:282:A:N3	1:CA:282:A:H2'	2.28	0.48
1:CA:345:C:H3'	28:DP:38:ARG:CZ	2.44	0.48
1:CA:370:C:H2'	1:CA:371:A:C8	2.49	0.48
1:CA:461:A:H3'	1:CA:462:G:C4'	2.43	0.48
1:CA:643:C:H2'	1:CA:644:U:H6	1.78	0.48
1:CA:821:G:H2'	1:CA:822:U:H6	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:903:G:H2'	1:CA:904:U:C6	2.49	0.48
1:CA:963:G:H2'	1:CA:964:A:C8	2.49	0.48
2:CC:26:LYS:CG	2:CC:27:GLU:N	2.75	0.48
2:CC:39:ARG:CZ	2:CC:56:ILE:HD11	2.43	0.48
3:CD:66:VAL:HG12	3:CD:70:GLN:OE1	2.13	0.48
8:CI:15:ALA:O	8:CI:66:VAL:HA	2.13	0.48
11:CL:106:VAL:HA	11:CL:107:LYS:HZ3	1.78	0.48
12:CM:21:ILE:HB	12:CM:24:VAL:CG2	2.31	0.48
21:CN:78:LEU:HD23	21:CN:82:LYS:O	2.13	0.48
14:CQ:7:LEU:O	14:CQ:60:ILE:HD13	2.14	0.48
16:CS:31:ARG:HD2	16:CS:31:ARG:N	2.28	0.48
34:D3:60:CYS:O	34:D3:61:LEU:HD23	2.14	0.48
22:DA:111:U:H2'	22:DA:112:G:H8	1.78	0.48
22:DA:64:G:H2'	22:DA:65:U:C6	2.49	0.48
23:DB:1010:A:H1'	23:DB:1153:C:O4'	2.13	0.48
23:DB:1631:G:H22	23:DB:1633:G:H3'	1.78	0.48
23:DB:1847:A:H4'	23:DB:1848:A:H8	1.78	0.48
23:DB:2047:C:H2'	23:DB:2048:G:H8	1.78	0.48
23:DB:2050:C:O2	26:DD:161:MET:HE1	2.14	0.48
23:DB:2259:U:H2'	23:DB:2260:C:C6	2.48	0.48
23:DB:267:C:H2'	23:DB:268:C:H6	1.79	0.48
23:DB:2742:G:O2'	23:DB:2743:U:H5'	2.14	0.48
23:DB:527:C:C5	23:DB:2779:U:H2'	2.48	0.48
23:DB:2895:G:H2'	23:DB:2896:C:C6	2.49	0.48
23:DB:438:G:H2'	23:DB:439:A:C8	2.46	0.48
23:DB:630:G:N2	23:DB:632:A:H3'	2.29	0.48
23:DB:838:C:O2'	23:DB:839:U:H5'	2.13	0.48
23:DB:873:C:H4'	38:DM:64:TRP:NE1	2.23	0.48
23:DB:921:C:H2'	23:DB:922:C:H6	1.78	0.48
23:DB:451:U:P	29:DE:47:LYS:HD2	2.53	0.48
29:DE:60:TRP:HE1	29:DE:71:GLY:HA2	1.79	0.48
47:DF:105:ILE:C	47:DF:108:PRO:HD2	2.33	0.48
47:DF:134:GLN:OE1	47:DF:136:ILE:HA	2.14	0.48
47:DF:135:ILE:C	47:DF:137:PHE:H	2.15	0.48
48:DG:129:GLU:C	48:DG:130:ILE:HG13	2.34	0.48
48:DG:122:ALA:CA	48:DG:132:LEU:HA	2.41	0.48
48:DG:43:LYS:HE2	48:DG:43:LYS:HA	1.94	0.48
40:DH:18:GLN:HE21	40:DH:39:ALA:CB	2.26	0.48
23:DB:1060:U:H5	24:DI:131:THR:HG22	1.78	0.48
24:DI:100:ILE:O	24:DI:139:VAL:HA	2.13	0.48
26:DD:21:SER:HB3	27:DK:72:PRO:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DK:60:ALA:CB	27:DK:86:LEU:HA	2.44	0.48
37:DL:89:VAL:HG13	37:DL:89:VAL:O	2.13	0.48
38:DM:90:GLU:HA	38:DM:90:GLU:OE1	2.12	0.48
42:DN:9:GLN:HA	42:DN:17:ARG:CD	2.44	0.48
43:DO:115:LEU:H	43:DO:115:LEU:CD2	2.27	0.48
28:DP:19:PHE:O	28:DP:20:ARG:HB2	2.14	0.48
28:DP:25:VAL:HB	28:DP:84:SER:O	2.14	0.48
44:DQ:86:SER:HB3	49:DR:51:VAL:HA	1.94	0.48
49:DR:49:ILE:HG12	49:DR:53:PHE:HA	1.94	0.48
50:DT:9:LYS:HB3	50:DT:9:LYS:HZ3	1.77	0.48
46:DU:6:ARG:O	46:DU:24:VAL:HB	2.13	0.48
52:DW:30:VAL:O	52:DW:30:VAL:HG13	2.13	0.48
1:AA:1048:G:H5''	21:AN:2:LYS:HD2	1.96	0.48
1:AA:1326:U:H2'	1:AA:1327:C:H6	1.79	0.48
1:AA:1414:U:H2'	1:AA:1415:G:H8	1.79	0.48
1:AA:382:A:H2'	1:AA:383:A:O4'	2.13	0.48
1:AA:502:A:H2'	1:AA:503:C:C6	2.49	0.48
1:AA:98:A:H2'	1:AA:99:C:H6	1.78	0.48
18:AB:14:HIS:CD2	18:AB:202:ASN:HB2	2.49	0.48
3:AD:151:GLN:O	3:AD:154:VAL:HG22	2.12	0.48
10:AK:41:LEU:HD22	10:AK:76:TYR:CD2	2.48	0.48
12:AM:10:ASP:C	12:AM:12:LYS:H	2.17	0.48
21:AN:6:LYS:HA	21:AN:62:ARG:HH11	1.78	0.48
13:AP:74:LEU:O	13:AP:78:VAL:HG12	2.12	0.48
14:AQ:18:LYS:H	14:AQ:50:ASN:ND2	2.12	0.48
17:AT:64:GLY:C	17:AT:66:ILE:H	2.16	0.48
33:B1:8:ILE:CD1	33:B1:9:LYS:H	2.27	0.48
22:BA:76:G:H1	22:BA:101:A:H61	1.62	0.48
23:BB:2487:G:H2'	23:BB:2488:G:C8	2.48	0.48
23:BB:308:G:O4'	23:BB:501:A:H5'	2.14	0.48
23:BB:30:G:OP1	44:BQ:4:LYS:HG2	2.14	0.48
23:BB:406:G:H2'	23:BB:407:G:O4'	2.13	0.48
23:BB:527:C:C5	23:BB:2779:U:H2'	2.47	0.48
23:BB:697:G:H2'	23:BB:698:C:C6	2.49	0.48
23:BB:789:A:H4'	23:BB:790:U:OP2	2.14	0.48
23:BB:831:G:O2'	23:BB:832:U:H5'	2.14	0.48
23:BB:958:U:O4	38:BM:16:ARG:HA	2.14	0.48
23:BB:968:C:H2'	23:BB:969:G:C8	2.48	0.48
25:BC:115:ILE:HB	25:BC:126:GLY:O	2.13	0.48
25:BC:152:GLN:C	25:BC:153:LEU:HD23	2.33	0.48
26:BD:32:ASN:HB3	26:BD:50:VAL:HG21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:113:VAL:C	29:BE:115:GLN:H	2.17	0.48
47:BF:168:LEU:O	47:BF:169:LEU:HB2	2.14	0.48
47:BF:87:LYS:C	47:BF:88:VAL:HG23	2.34	0.48
41:BJ:37:ARG:HE	41:BJ:110:PRO:HG3	1.79	0.48
27:BK:21:CYS:HB2	27:BK:39:ILE:HG21	1.96	0.48
42:BN:17:ARG:C	42:BN:19:ALA:H	2.17	0.48
52:BW:61:LYS:HB3	52:BW:62:ALA:H	1.48	0.48
30:BY:35:VAL:HG22	30:BY:36:GLU:N	2.28	0.48
30:BY:50:VAL:HB	30:BY:53:MET:HB2	1.94	0.48
23:BB:200:U:H4'	51:BZ:22:LEU:HB2	1.96	0.48
51:BZ:59:ILE:CD1	51:BZ:67:VAL:HG21	2.42	0.48
51:BZ:5:CYS:SG	51:BZ:8:THR:HG23	2.53	0.48
1:CA:1020:G:H2'	1:CA:1021:A:H5'	1.95	0.48
1:CA:1387:G:O2'	1:CA:1388:C:H5'	2.13	0.48
1:CA:211:G:H3'	1:CA:211:G:N3	2.29	0.48
1:CA:254:G:H4'	14:CQ:19:SER:OG	2.13	0.48
1:CA:56:U:H2'	1:CA:57:G:C8	2.49	0.48
2:CC:180:ASP:OD2	2:CC:203:LYS:HB2	2.14	0.48
2:CC:51:VAL:HG21	2:CC:67:ILE:HG23	1.94	0.48
3:CD:35:GLN:O	3:CD:37:PRO:HD3	2.13	0.48
1:CA:8:A:H5''	4:CE:125:LYS:HB3	1.94	0.48
4:CE:76:ASN:HB2	4:CE:81:GLN:NE2	2.29	0.48
5:CF:9:MET:HA	5:CF:58:HIS:O	2.13	0.48
9:CJ:52:LEU:HG	9:CJ:62:ARG:NE	2.19	0.48
13:CP:54:LEU:HD22	13:CP:80:LYS:HE3	1.96	0.48
15:CR:20:ILE:HG22	15:CR:53:GLN:CD	2.33	0.48
17:CT:54:GLN:N	17:CT:55:PRO:HD2	2.28	0.48
34:D3:50:SER:C	34:D3:52:GLY:H	2.15	0.48
22:DA:29:A:OP1	43:DO:31:THR:HB	2.13	0.48
23:DB:1571:A:H2'	23:DB:1572:A:H8	1.78	0.48
23:DB:1747:U:H2'	23:DB:1748:C:H6	1.77	0.48
23:DB:1805:A:O2'	23:DB:1806:C:H5'	2.14	0.48
23:DB:2538:C:H2'	23:DB:2539:C:H6	1.79	0.48
23:DB:2544:G:O2'	23:DB:2545:G:H5'	2.13	0.48
23:DB:499:U:H2'	23:DB:500:G:O4'	2.14	0.48
23:DB:811:U:O2	23:DB:1250:G:H2'	2.13	0.48
25:DC:130:PRO:HA	25:DC:187:CYS:O	2.14	0.48
23:DB:1820:U:C4	25:DC:158:GLY:HA3	2.48	0.48
26:DD:37:VAL:CG2	26:DD:91:THR:HA	2.42	0.48
40:DH:135:HIS:CD2	40:DH:137:GLU:H	2.32	0.48
24:DI:129:GLU:HB3	24:DI:133:ARG:NH1	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DJ:7:LYS:C	41:DJ:9:GLU:H	2.17	0.48
28:DP:89:GLY:H	28:DP:112:ARG:NH1	2.11	0.48
46:DU:34:ILE:HG12	46:DU:63:ALA:CB	2.43	0.48
1:AA:1107:C:C4	1:AA:1108:G:N7	2.82	0.48
1:AA:1198:G:H2'	1:AA:1199:U:C6	2.49	0.48
1:AA:1360:A:H61	1:AA:1361:G:N2	2.11	0.48
1:AA:590:U:H2'	1:AA:591:U:C6	2.48	0.48
1:AA:647:C:H2'	1:AA:648:A:H8	1.79	0.48
1:AA:68:G:H2'	1:AA:69:G:O4'	2.14	0.48
1:AA:947:G:H2'	1:AA:948:C:H6	1.76	0.48
18:AB:110:ILE:HG12	18:AB:147:LEU:HD13	1.95	0.48
3:AD:110:ARG:HH11	3:AD:110:ARG:HG3	1.79	0.48
3:AD:90:LEU:N	3:AD:90:LEU:HD22	2.28	0.48
6:AG:132:THR:O	6:AG:135:LYS:HB3	2.14	0.48
7:AH:94:VAL:HG21	7:AH:100:ILE:O	2.13	0.48
8:AI:97:LEU:O	8:AI:102:PHE:HB2	2.13	0.48
20:AO:56:LEU:O	20:AO:60:VAL:HG23	2.14	0.48
11:AL:7:VAL:HG22	14:AQ:33:TYR:HD1	1.79	0.48
36:B2:10:LEU:HD21	36:B2:14:ARG:CZ	2.43	0.48
34:B3:49:VAL:HG11	34:B3:54:LEU:CD1	2.42	0.48
23:BB:1080:A:O2'	23:BB:1081:U:H5'	2.13	0.48
23:BB:1144:A:O2'	23:BB:1145:C:H5'	2.14	0.48
23:BB:1259:G:H2'	23:BB:1260:A:C8	2.49	0.48
23:BB:1530:G:N3	23:BB:1530:G:H2'	2.29	0.48
23:BB:1636:U:H2'	23:BB:1637:A:C8	2.48	0.48
23:BB:170:U:H2'	23:BB:171:U:C6	2.49	0.48
23:BB:1797:G:O2'	23:BB:1798:U:H5'	2.13	0.48
23:BB:182:A:O2'	23:BB:183:C:H5'	2.14	0.48
23:BB:2037:A:H2'	23:BB:2038:G:C8	2.49	0.48
23:BB:2041:U:H2'	23:BB:2042:A:H8	1.78	0.48
23:BB:2428:G:H5''	23:BB:2429:G:OP1	2.14	0.48
23:BB:796:C:H2'	23:BB:797:G:H8	1.78	0.48
23:BB:90:U:H2'	23:BB:91:A:C2	2.49	0.48
23:BB:992:C:H2'	23:BB:993:G:H8	1.78	0.48
25:BC:183:VAL:HG22	25:BC:184:GLU:N	2.28	0.48
25:BC:199:HIS:C	25:BC:201:LEU:H	2.17	0.48
26:BD:97:SER:OG	26:BD:98:VAL:N	2.47	0.48
29:BE:3:LEU:CB	29:BE:12:LEU:HB3	2.44	0.48
29:BE:28:VAL:HG23	29:BE:29:HIS:H	1.78	0.48
29:BE:60:TRP:C	29:BE:62:GLN:H	2.15	0.48
47:BF:76:PHE:HD2	47:BF:78:ILE:HD13	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BF:90:LEU:C	47:BF:91:ARG:HD3	2.34	0.48
47:BF:3:LEU:HD12	47:BF:96:TRP:CD1	2.49	0.48
40:BH:44:ILE:HA	40:BH:51:ARG:NH2	2.27	0.48
24:BI:52:LEU:HD12	24:BI:52:LEU:N	2.29	0.48
27:BK:19:VAL:HB	27:BK:41:ILE:CD1	2.43	0.48
27:BK:92:GLU:HB2	27:BK:93:GLN:HE21	1.78	0.48
37:BL:122:VAL:HB	37:BL:143:GLU:OE2	2.14	0.48
44:BQ:105:PHE:O	44:BQ:109:VAL:HG23	2.13	0.48
46:BU:95:PHE:HB2	46:BU:100:GLU:HB3	1.95	0.48
52:BW:48:ALA:O	52:BW:61:LYS:HB2	2.13	0.48
1:CA:1102:A:H2'	1:CA:1103:C:C6	2.49	0.48
1:CA:1213:A:H2'	1:CA:1215:G:N7	2.29	0.48
1:CA:1253:G:C6	1:CA:1285:A:N6	2.81	0.48
1:CA:1288:A:H2'	1:CA:1289:A:O4'	2.13	0.48
1:CA:284:C:H2'	1:CA:285:C:H6	1.79	0.48
1:CA:284:C:O2'	1:CA:285:C:H5'	2.14	0.48
1:CA:309:A:O2'	1:CA:310:G:H5'	2.13	0.48
1:CA:376:G:H2'	1:CA:377:G:H8	1.78	0.48
1:CA:438:U:H4'	3:CD:119:HIS:HD2	1.78	0.48
1:CA:601:G:H2'	1:CA:602:A:H8	1.78	0.48
1:CA:634:C:H2'	1:CA:635:A:C8	2.49	0.48
18:CB:127:LYS:HD2	18:CB:128:LEU:HB2	1.95	0.48
18:CB:131:LYS:O	18:CB:134:LEU:HG	2.13	0.48
5:CF:21:MET:HB3	5:CF:25:TYR:CE1	2.49	0.48
8:CI:114:LYS:HD3	8:CI:114:LYS:N	2.29	0.48
8:CI:6:TYR:CG	8:CI:7:GLY:N	2.81	0.48
20:CO:28:GLN:O	20:CO:32:LEU:HD23	2.12	0.48
17:CT:19:HIS:O	17:CT:23:ARG:HG2	2.14	0.48
10:CK:110:THR:CG2	19:CU:4:LYS:HA	2.44	0.48
34:D3:60:CYS:C	34:D3:62:PRO:HD3	2.34	0.48
23:DB:132:G:H2'	23:DB:133:U:H6	1.78	0.48
23:DB:1477:A:H2'	23:DB:1478:G:O4'	2.13	0.48
23:DB:1818:U:C4	25:DC:152:GLN:HB3	2.49	0.48
23:DB:2105:U:H2'	23:DB:2106:U:H6	1.79	0.48
23:DB:464:U:H2'	23:DB:465:G:O4'	2.14	0.48
23:DB:1205:A:N6	29:DE:165:HIS:HB2	2.15	0.48
29:DE:28:VAL:HG23	29:DE:29:HIS:H	1.78	0.48
29:DE:28:VAL:HG23	29:DE:29:HIS:N	2.28	0.48
29:DE:62:GLN:CG	29:DE:63:LYS:H	2.27	0.48
47:DF:148:VAL:O	47:DF:149:ARG:HG2	2.13	0.48
23:DB:1080:A:H4'	24:DI:126:ARG:HD2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DJ:106:LYS:HE3	41:DJ:106:LYS:O	2.13	0.48
27:DK:60:ALA:HA	27:DK:87:LEU:HD22	1.96	0.48
37:DL:57:LEU:O	37:DL:61:LEU:HD13	2.13	0.48
42:DN:79:LEU:O	42:DN:80:PHE:HB2	2.12	0.48
42:DN:90:ARG:HG2	42:DN:94:TYR:HD1	1.78	0.48
43:DO:30:ARG:HG2	43:DO:31:THR:N	2.29	0.48
44:DQ:105:PHE:O	44:DQ:109:VAL:HG23	2.14	0.48
46:DU:40:LEU:HA	46:DU:61:GLU:HA	1.95	0.48
46:DU:3:LYS:O	46:DU:4:ILE:HD13	2.14	0.48
52:DW:33:GLY:O	52:DW:34:SER:HB2	2.13	0.48
30:DY:11:SER:OG	30:DY:13:ILE:HG13	2.13	0.48
1:AA:1047:G:H21	1:AA:1215:G:C5'	2.25	0.48
1:AA:1207:G:H2'	1:AA:1208:C:C6	2.49	0.48
1:AA:1308:U:H5''	12:AM:96:VAL:HG23	1.95	0.48
1:AA:221:C:O2'	1:AA:222:C:H5'	2.14	0.48
1:AA:376:G:H2'	1:AA:377:G:H8	1.79	0.48
1:AA:453:G:H2'	1:AA:454:G:C8	2.48	0.48
1:AA:488:C:O2'	1:AA:489:C:H5'	2.14	0.48
1:AA:594:U:H2'	1:AA:595:A:O4'	2.14	0.48
18:AB:120:SER:CA	18:AB:125:PHE:HB3	2.43	0.48
3:AD:129:VAL:HG12	3:AD:130:ASN:N	2.29	0.48
3:AD:66:VAL:HG12	3:AD:70:GLN:OE1	2.13	0.48
4:AE:81:GLN:HG2	4:AE:148:SER:HA	1.96	0.48
5:AF:21:MET:HB3	5:AF:25:TYR:CE1	2.49	0.48
1:AA:1232:U:OP1	8:AI:127:SER:HB3	2.14	0.48
12:AM:15:VAL:HG23	12:AM:33:LEU:HG	1.94	0.48
12:AM:89:ARG:HD3	12:AM:94:LEU:HB2	1.95	0.48
17:AT:38:ILE:HD11	17:AT:82:ILE:HA	1.94	0.48
33:B1:16:THR:HG21	33:B1:39:ASP:CG	2.34	0.48
22:BA:22:U:H2'	22:BA:23:G:C8	2.49	0.48
23:BB:1283:G:N2	23:BB:1285:A:H3'	2.29	0.48
23:BB:1469:A:H2'	23:BB:1470:A:C8	2.49	0.48
23:BB:2311:A:H3'	23:BB:2312:U:C6	2.48	0.48
23:BB:2649:C:H2'	23:BB:2650:U:C6	2.49	0.48
23:BB:346:A:H3'	23:BB:347:A:H8	1.79	0.48
23:BB:557:C:H2'	23:BB:558:U:C6	2.49	0.48
25:BC:163:ILE:HG22	25:BC:164:VAL:N	2.29	0.48
23:BB:1655:A:H5'	26:BD:118:PHE:HB2	1.96	0.48
26:BD:138:LEU:HD22	26:BD:138:LEU:N	2.29	0.48
26:BD:107:VAL:H	26:BD:206:ALA:H	1.62	0.48
26:BD:29:VAL:O	26:BD:185:ASN:HB3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BF:52:ALA:O	47:BF:56:LEU:HB2	2.14	0.48
23:BB:2529:G:H5'	48:BG:174:LYS:HB2	1.95	0.48
40:BH:99:ILE:HD13	40:BH:144:VAL:HB	1.95	0.48
24:BI:124:MET:O	24:BI:128:ILE:HG12	2.14	0.48
41:BJ:13:ARG:HB3	41:BJ:53:TYR:CD2	2.49	0.48
27:BK:105:ARG:HD3	27:BK:105:ARG:N	2.21	0.48
27:BK:109:SER:OG	27:BK:111:LYS:HG2	2.13	0.48
27:BK:73:ASP:OD2	27:BK:75:SER:HB3	2.14	0.48
37:BL:109:LYS:HB2	37:BL:111:ILE:HD13	1.96	0.48
44:BQ:56:PHE:O	44:BQ:59:LEU:HB3	2.13	0.48
45:BS:45:VAL:HG23	45:BS:46:LEU:HD23	1.95	0.48
35:BV:59:GLU:HG2	35:BV:60:VAL:N	2.28	0.48
23:BB:988:A:P	30:BY:11:SER:HB3	2.54	0.48
51:BZ:3:ARG:HG3	51:BZ:50:ARG:HH11	1.77	0.48
1:CA:1045:C:O2'	1:CA:1046:A:H5'	2.14	0.48
1:CA:1119:C:O2'	1:CA:1120:C:H5'	2.14	0.48
1:CA:1306:A:N6	1:CA:1331:G:H1'	2.29	0.48
1:CA:1368:A:O2'	1:CA:1369:C:H5'	2.14	0.48
1:CA:777:A:H2'	1:CA:778:G:C8	2.49	0.48
18:CB:86:CYS:SG	18:CB:87:ASP:N	2.87	0.48
2:CC:152:VAL:HB	2:CC:165:GLU:CB	2.44	0.48
2:CC:21:TRP:CD2	2:CC:58:ARG:HD3	2.49	0.48
2:CC:46:LEU:N	2:CC:46:LEU:HD22	2.28	0.48
7:CH:99:GLY:HA2	7:CH:129:ALA:HA	1.96	0.48
21:CN:73:LEU:HD12	21:CN:83:VAL:HG21	1.95	0.48
2:CC:18:ASN:HB2	21:CN:90:GLY:O	2.14	0.48
20:CO:33:THR:HG23	20:CO:63:ARG:NH1	2.29	0.48
34:D3:28:LEU:O	34:D3:29:ARG:HB3	2.14	0.48
32:D4:15:LYS:O	32:D4:16:ILE:HB	2.14	0.48
23:DB:2014:A:H2'	23:DB:2015:A:C8	2.49	0.48
23:DB:2348:U:OP1	34:D3:37:THR:HG21	2.14	0.48
23:DB:2675:A:N1	23:DB:2732:G:O6	2.47	0.48
23:DB:2773:C:O2'	23:DB:2774:C:H5'	2.14	0.48
25:DC:68:ARG:HH21	25:DC:190:THR:CG2	2.27	0.48
23:DB:600:G:C1'	29:DE:100:MET:HG2	2.39	0.48
47:DF:71:LYS:O	47:DF:72:SER:HB3	2.14	0.48
40:DH:62:LEU:HG	40:DH:66:ASN:HD21	1.77	0.48
27:DK:68:GLY:HA3	27:DK:78:ARG:HB3	1.96	0.48
43:DO:111:ARG:HD2	43:DO:117:PHE:CD1	2.48	0.48
28:DP:3:ILE:HD13	28:DP:4:ILE:N	2.29	0.48
44:DQ:94:LEU:C	44:DQ:96:ASP:N	2.66	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:DT:8:LEU:HD22	50:DT:46:ALA:HA	1.95	0.48
39:DX:45:GLN:O	39:DX:47:ARG:N	2.46	0.48
1:AA:1478:U:H2'	1:AA:1479:C:C6	2.49	0.48
1:AA:1515:G:O2'	1:AA:1516:G:H5'	2.13	0.48
1:AA:299:G:H2'	1:AA:300:A:C8	2.48	0.48
1:AA:643:C:H2'	1:AA:644:U:H6	1.79	0.48
1:AA:846:G:H2'	1:AA:846:G:N3	2.29	0.48
18:AB:102:ASN:HD21	18:AB:105:THR:HB	1.77	0.48
18:AB:183:PHE:CD1	18:AB:183:PHE:N	2.82	0.48
4:AE:146:MET:HG3	4:AE:146:MET:O	2.12	0.48
6:AG:94:ARG:HH11	6:AG:98:LEU:HD21	1.77	0.48
12:AM:12:LYS:O	12:AM:43:LYS:HA	2.14	0.48
16:AS:31:ARG:HB3	16:AS:33:TRP:CZ2	2.49	0.48
36:B2:31:LEU:CD2	36:B2:42:LEU:HD12	2.44	0.48
23:BB:1409:U:H2'	23:BB:1410:G:H8	1.79	0.48
23:BB:143:C:H6	23:BB:143:C:O5'	1.97	0.48
23:BB:1607:C:H5''	23:BB:1608:A:H5'	1.96	0.48
23:BB:163:C:H2'	23:BB:164:C:O4'	2.14	0.48
23:BB:1779:U:C5	23:BB:1784:A:N7	2.82	0.48
23:BB:2305:U:H2'	23:BB:2306:C:C6	2.48	0.48
23:BB:2305:U:H5''	47:BF:130:GLY:CA	2.43	0.48
23:BB:2899:A:H2'	23:BB:2900:A:O4'	2.14	0.48
23:BB:702:U:H2'	23:BB:703:U:C6	2.49	0.48
23:BB:973:A:OP1	23:BB:973:A:H8	1.96	0.48
25:BC:270:ARG:HB3	25:BC:270:ARG:HH11	1.79	0.48
25:BC:34:GLU:O	25:BC:34:GLU:HG3	2.13	0.48
25:BC:42:ARG:CZ	25:BC:48:ILE:HD11	2.43	0.48
29:BE:88:ARG:O	29:BE:90:GLN:HG3	2.14	0.48
47:BF:102:LEU:HD13	47:BF:103:ILE:N	2.29	0.48
47:BF:134:GLN:OE1	47:BF:136:ILE:HA	2.13	0.48
47:BF:64:PRO:HA	47:BF:88:VAL:CG1	2.43	0.48
48:BG:59:ASP:O	48:BG:63:GLN:HB2	2.14	0.48
41:BJ:72:LYS:HB2	41:BJ:89:PHE:H	1.77	0.48
37:BL:4:ASN:N	37:BL:4:ASN:HD22	2.11	0.48
43:BO:66:GLY:O	43:BO:102:ARG:HD3	2.14	0.48
44:BQ:24:TYR:CD1	44:BQ:25:GLY:N	2.82	0.48
50:BT:40:LYS:HA	50:BT:43:ILE:HG22	1.96	0.48
1:CA:299:G:C6	1:CA:300:A:C6	3.02	0.48
18:CB:221:ARG:HH11	18:CB:221:ARG:HB3	1.79	0.48
18:CB:30:ILE:HG23	18:CB:39:ILE:O	2.14	0.48
6:CG:31:VAL:HG22	6:CG:32:ASP:CG	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:37:ARG:NH1	9:CJ:37:ARG:HA	2.28	0.48
10:CK:30:ILE:HG22	10:CK:45:THR:HB	1.95	0.48
11:CL:13:ARG:HB2	11:CL:14:LYS:H	1.50	0.48
21:CN:14:ALA:O	21:CN:18:LYS:HG3	2.13	0.48
19:CU:16:ARG:HH12	19:CU:19:LYS:NZ	2.11	0.48
34:D3:2:LYS:H	34:D3:2:LYS:HG2	1.44	0.48
22:DA:22:U:H2'	22:DA:23:G:C8	2.49	0.48
23:DB:1005:C:H2'	23:DB:1006:C:C6	2.49	0.48
23:DB:1099:G:H3'	24:DI:2:LYS:HA	1.94	0.48
23:DB:1290:C:O2'	23:DB:1291:C:H5'	2.13	0.48
23:DB:1866:A:H2'	23:DB:1867:G:O4'	2.13	0.48
23:DB:1871:A:H2'	23:DB:1872:A:H8	1.77	0.48
23:DB:970:U:H2'	23:DB:971:G:H8	1.76	0.48
25:DC:180:MET:HB2	25:DC:268:ARG:HB3	1.96	0.48
25:DC:79:ARG:HD2	25:DC:81:GLU:OE2	2.14	0.48
26:DD:183:GLU:OE1	26:DD:183:GLU:N	2.47	0.48
26:DD:47:ALA:HB2	26:DD:83:ARG:HD2	1.95	0.48
29:DE:113:VAL:C	29:DE:115:GLN:H	2.16	0.48
47:DF:49:LEU:HD21	47:DF:84:ILE:O	2.13	0.48
48:DG:106:LEU:C	48:DG:108:PHE:H	2.17	0.48
40:DH:4:ILE:O	40:DH:36:ALA:HA	2.13	0.48
24:DI:126:ARG:NH1	24:DI:126:ARG:HB3	2.28	0.48
24:DI:2:LYS:N	24:DI:2:LYS:HD2	2.29	0.48
41:DJ:72:LYS:HB2	41:DJ:89:PHE:H	1.79	0.48
27:DK:25:LEU:CD2	27:DK:40:LYS:HB2	2.37	0.48
38:DM:102:LEU:N	38:DM:102:LEU:HD22	2.28	0.48
22:DA:113:C:H1'	43:DO:46:GLU:HA	1.96	0.48
44:DQ:93:ILE:O	44:DQ:96:ASP:HB3	2.14	0.48
49:DR:22:LEU:N	49:DR:22:LEU:HD23	2.29	0.48
1:AA:174:A:O2'	1:AA:175:C:H5'	2.14	0.47
1:AA:857:C:H2'	1:AA:858:G:O4'	2.14	0.47
1:AA:933:G:N2	1:AA:935:A:O4'	2.47	0.47
18:AB:13:VAL:O	18:AB:14:HIS:CB	2.62	0.47
3:AD:96:ARG:HH12	3:AD:133:SER:HA	1.76	0.47
4:AE:158:LYS:HZ3	7:AH:63:LYS:HD3	1.78	0.47
6:AG:106:ALA:HB1	6:AG:132:THR:OG1	2.13	0.47
8:AI:90:ASP:O	8:AI:93:LEU:HG	2.13	0.47
9:AJ:80:THR:O	9:AJ:84:VAL:HG23	2.14	0.47
12:AM:47:LEU:CG	12:AM:52:ILE:HD13	2.29	0.47
23:BB:1173:U:H2'	23:BB:1174:U:H5'	1.96	0.47
23:BB:1234:U:H2'	23:BB:1235:G:O4'	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1365:A:N3	23:BB:1365:A:H2'	2.28	0.47
23:BB:2721:A:H2'	23:BB:2722:G:C8	2.49	0.47
23:BB:499:U:H2'	23:BB:500:G:O4'	2.14	0.47
23:BB:553:G:C2'	23:BB:554:U:H5'	2.44	0.47
25:BC:141:HIS:NE2	25:BC:194:VAL:HA	2.29	0.47
47:BF:41:GLU:CG	47:BF:48:LEU:HD11	2.44	0.47
48:BG:154:GLU:C	48:BG:156:TYR:H	2.18	0.47
48:BG:26:LYS:HB2	48:BG:32:LEU:HG	1.96	0.47
41:BJ:117:ALA:O	41:BJ:120:ARG:HG2	2.14	0.47
41:BJ:14:ASP:O	41:BJ:53:TYR:HB2	2.14	0.47
37:BL:116:VAL:HG22	37:BL:117:THR:N	2.30	0.47
38:BM:101:VAL:HG22	38:BM:101:VAL:O	2.13	0.47
38:BM:34:LYS:O	38:BM:128:THR:HG22	2.13	0.47
42:BN:59:SER:O	42:BN:63:ARG:HB2	2.13	0.47
28:BP:19:PHE:CE2	28:BP:25:VAL:HG11	2.49	0.47
44:BQ:79:ILE:HD13	44:BQ:79:ILE:C	2.35	0.47
44:BQ:93:ILE:HG23	44:BQ:94:LEU:N	2.29	0.47
49:BR:58:VAL:HG22	49:BR:59:ILE:N	2.24	0.47
49:BR:14:VAL:HG21	49:BR:98:ILE:HD13	1.96	0.47
46:BU:82:VAL:N	46:BU:96:LYS:HZ2	2.06	0.47
35:BV:75:GLN:HB3	35:BV:76:ASP:H	1.43	0.47
30:BY:35:VAL:HG11	30:BY:37:ARG:HH12	1.79	0.47
1:CA:1219:A:H2'	1:CA:1220:G:H8	1.78	0.47
1:CA:1241:G:H2'	1:CA:1242:G:C8	2.41	0.47
1:CA:131:A:H2'	1:CA:132:C:C6	2.49	0.47
1:CA:238:A:C2'	1:CA:239:U:H5''	2.43	0.47
1:CA:424:G:O2'	1:CA:425:G:H5'	2.14	0.47
1:CA:738:C:H2'	1:CA:739:C:H6	1.79	0.47
18:CB:48:MET:HA	18:CB:51:GLU:OE2	2.14	0.47
18:CB:80:LYS:HE2	18:CB:81:ASP:OD1	2.13	0.47
2:CC:107:LYS:C	2:CC:109:GLU:H	2.17	0.47
2:CC:111:ASP:O	2:CC:115:VAL:HG23	2.14	0.47
2:CC:139:ASN:N	2:CC:139:ASN:HD22	2.12	0.47
2:CC:188:ALA:O	2:CC:194:VAL:HA	2.14	0.47
5:CF:93:LYS:O	5:CF:94:HIS:HB2	2.14	0.47
20:CO:40:GLN:O	20:CO:44:ALA:N	2.45	0.47
31:D0:27:LEU:HD13	31:D0:37:HIS:O	2.13	0.47
31:D0:2:VAL:HG12	31:D0:3:GLN:N	2.29	0.47
36:D2:44:VAL:O	36:D2:45:SER:C	2.51	0.47
34:D3:18:LYS:HD2	34:D3:19:GLY:H	1.78	0.47
23:DB:1017:G:O2'	23:DB:1018:U:H5'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:118:A:P	23:DB:119:A:H5''	2.54	0.47
23:DB:1287:A:H3'	23:DB:1288:G:H21	1.76	0.47
23:DB:170:U:H2'	23:DB:171:U:C6	2.48	0.47
23:DB:1133:A:N6	23:DB:2025:C:O2'	2.46	0.47
23:DB:2300:C:H2'	23:DB:2301:C:H6	1.79	0.47
23:DB:2365:G:O2'	52:DW:59:PHE:HE1	1.97	0.47
23:DB:2514:U:H5''	41:DJ:81:ILE:HD12	1.96	0.47
23:DB:433:C:O2'	23:DB:434:U:H5'	2.14	0.47
26:DD:107:VAL:H	26:DD:206:ALA:H	1.62	0.47
26:DD:34:VAL:HA	26:DD:50:VAL:HG12	1.95	0.47
26:DD:63:PRO:C	26:DD:65:ALA:H	2.18	0.47
29:DE:126:VAL:HG13	29:DE:127:GLU:N	2.29	0.47
29:DE:67:ARG:HG2	29:DE:67:ARG:NH1	2.26	0.47
47:DF:139:GLU:CG	47:DF:140:ILE:H	2.26	0.47
47:DF:41:GLU:CG	47:DF:48:LEU:HD11	2.44	0.47
48:DG:84:LYS:HB2	48:DG:132:LEU:N	2.29	0.47
38:DM:21:ALA:HB1	38:DM:100:LYS:HE2	1.95	0.47
38:DM:73:ILE:HG21	38:DM:91:TYR:CZ	2.49	0.47
43:DO:14:ALA:C	43:DO:16:ARG:H	2.17	0.47
45:DS:37:THR:HG22	45:DS:48:LYS:HE3	1.96	0.47
50:DT:50:LEU:O	50:DT:51:PHE:HB2	2.14	0.47
46:DU:73:ASN:C	46:DU:75:ALA:H	2.17	0.47
23:DB:2331:G:O4'	52:DW:39:GLN:HA	2.14	0.47
52:DW:39:GLN:HG2	52:DW:40:ARG:H	1.77	0.47
52:DW:9:THR:HG23	52:DW:10:ARG:CD	2.44	0.47
52:DW:9:THR:OG1	52:DW:10:ARG:N	2.44	0.47
30:DY:6:ILE:H	30:DY:6:ILE:CD1	2.24	0.47
51:DZ:15:GLY:O	51:DZ:27:ARG:HG3	2.14	0.47
51:DZ:27:ARG:O	51:DZ:28:ARG:CB	2.62	0.47
51:DZ:33:LEU:H	51:DZ:52:SER:HB3	1.78	0.47
1:AA:1038:C:H2'	1:AA:1039:G:C8	2.49	0.47
1:AA:1062:U:H2'	1:AA:1063:C:C6	2.49	0.47
1:AA:1085:U:H3'	1:AA:1086:U:H5	1.78	0.47
1:AA:1414:U:O2'	1:AA:1415:G:H5'	2.14	0.47
1:AA:1434:A:H2'	1:AA:1435:G:O4'	2.14	0.47
1:AA:1519:A:H3'	1:AA:1520:C:C5'	2.44	0.47
1:AA:781:A:O2'	1:AA:1522:U:O2	2.30	0.47
1:AA:160:A:H2'	1:AA:161:A:O4'	2.14	0.47
1:AA:204:G:N2	1:AA:466:A:N6	2.62	0.47
1:AA:230:G:H5''	13:AP:31:ARG:NH2	2.29	0.47
1:AA:501:C:H1'	1:AA:549:C:H1'	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:559:A:H4'	1:AA:560:A:H3'	1.97	0.47
1:AA:664:G:N2	1:AA:741:G:H1	2.01	0.47
1:AA:767:A:H2'	1:AA:768:A:C8	2.49	0.47
1:AA:923:A:H2'	1:AA:924:C:H6	1.79	0.47
18:AB:95:TRP:CH2	18:AB:171:ALA:HA	2.49	0.47
18:AB:162:VAL:HG13	18:AB:184:ALA:HB2	1.96	0.47
18:AB:40:ILE:O	18:AB:42:LEU:N	2.47	0.47
2:AC:37:LYS:O	2:AC:41:TYR:HB2	2.13	0.47
4:AE:149:PRO:HG2	4:AE:150:GLU:H	1.79	0.47
8:AI:5:TYR:HB2	8:AI:20:ILE:HB	1.96	0.47
9:AJ:70:HIS:HB3	9:AJ:72:ARG:HH21	1.79	0.47
1:AA:568:G:O6	11:AL:1:ALA:HB2	2.14	0.47
11:AL:41:PRO:HA	11:AL:88:ASP:O	2.13	0.47
16:AS:4:LEU:O	16:AS:5:LYS:C	2.52	0.47
17:AT:25:SER:O	17:AT:29:THR:HG23	2.14	0.47
22:BA:64:G:O2'	22:BA:65:U:H5'	2.14	0.47
23:BB:1431:A:H2'	23:BB:1432:G:H8	1.78	0.47
23:BB:1550:C:H2'	23:BB:1551:A:C8	2.50	0.47
23:BB:1439:A:C6	23:BB:1552:A:N7	2.82	0.47
23:BB:1672:A:C6	23:BB:1673:G:C6	3.02	0.47
23:BB:2135:A:H61	23:BB:2156:G:C2'	2.27	0.47
23:BB:2254:C:H2'	23:BB:2255:G:O4'	2.13	0.47
23:BB:2350:C:H2'	23:BB:2351:G:C8	2.49	0.47
23:BB:242:G:H5''	34:B3:63:TYR:CD2	2.49	0.47
23:BB:2617:U:C2'	23:BB:2618:G:H5'	2.44	0.47
23:BB:2831:G:H1'	23:BB:2883:A:C2	2.48	0.47
23:BB:704:G:H1'	23:BB:727:A:H61	1.79	0.47
29:BE:15:SER:HB3	29:BE:18:THR:OG1	2.15	0.47
47:BF:102:LEU:O	47:BF:103:ILE:HB	2.13	0.47
48:BG:8:VAL:CG1	48:BG:49:LEU:HB3	2.39	0.47
40:BH:73:ASN:HB2	40:BH:141:LYS:HE2	1.96	0.47
40:BH:94:ILE:HG21	40:BH:144:VAL:HG11	1.94	0.47
23:BB:1250:G:OP2	37:BL:21:ARG:NH2	2.47	0.47
42:BN:96:ARG:NE	42:BN:116:VAL:HA	2.28	0.47
43:BO:7:ARG:HG3	43:BO:96:GLY:C	2.34	0.47
28:BP:12:MET:HG2	28:BP:54:LEU:HA	1.96	0.47
44:BQ:23:TYR:N	44:BQ:23:TYR:CD2	2.81	0.47
1:CA:1084:G:H5'	1:CA:1102:A:OP2	2.14	0.47
1:CA:410:G:O5'	1:CA:410:G:H8	1.97	0.47
1:CA:921:U:O2	4:CE:23:THR:HG23	2.14	0.47
2:CC:149:LYS:HA	2:CC:168:ARG:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:CG:144:ALA:O	6:CG:145:GLU:HB3	2.14	0.47
6:CG:144:ALA:O	6:CG:146:ALA:N	2.46	0.47
7:CH:101:ALA:HB3	7:CH:112:ASP:HB3	1.95	0.47
21:CN:98:ALA:HB1	21:CN:100:TRP:CZ3	2.48	0.47
20:CO:43:PHE:CD1	20:CO:56:LEU:HD22	2.49	0.47
14:CQ:26:ARG:HG2	14:CQ:39:ARG:O	2.14	0.47
17:CT:49:ALA:HA	17:CT:52:GLU:HB3	1.96	0.47
36:D2:9:VAL:CG1	36:D2:10:LEU:H	2.27	0.47
22:DA:75:G:H2'	22:DA:76:G:C8	2.48	0.47
23:DB:111:A:O2'	23:DB:112:U:H5'	2.14	0.47
23:DB:1521:G:H3'	23:DB:1522:A:H2'	1.96	0.47
23:DB:1559:U:H3'	23:DB:1560:G:H5'	1.96	0.47
23:DB:1647:U:H3'	23:DB:1647:U:OP1	2.14	0.47
23:DB:163:C:H2'	23:DB:164:C:O4'	2.15	0.47
23:DB:170:U:O2'	23:DB:171:U:H5'	2.14	0.47
23:DB:1716:U:H2'	23:DB:1717:A:H8	1.79	0.47
1:CA:1484:C:O2'	23:DB:1961:C:H5'	2.14	0.47
23:DB:2104:C:H2'	23:DB:2105:U:C6	2.49	0.47
23:DB:2559:C:O2'	23:DB:2560:A:H5'	2.13	0.47
23:DB:256:A:H2'	23:DB:257:C:H6	1.79	0.47
23:DB:2655:G:H1'	23:DB:2656:U:H5	1.79	0.47
23:DB:2877:G:O2'	23:DB:2878:U:H5'	2.14	0.47
23:DB:2876:G:H2'	23:DB:2877:G:O4'	2.15	0.47
23:DB:406:G:H2'	23:DB:407:G:O4'	2.13	0.47
23:DB:651:G:O5'	34:D3:17:GLY:HA3	2.14	0.47
23:DB:833:A:H2'	23:DB:834:G:C8	2.49	0.47
23:DB:864:G:O2'	23:DB:865:C:H5'	2.15	0.47
23:DB:1820:U:OP1	25:DC:176:ARG:HD2	2.14	0.47
47:DF:141:ASP:CB	47:DF:144:LYS:HB2	2.42	0.47
48:DG:108:PHE:HE1	48:DG:151:ARG:HH11	1.62	0.47
24:DI:17:ALA:O	24:DI:18:ASN:HB3	2.13	0.47
27:DK:71:ARG:HB3	27:DK:72:PRO:CD	2.38	0.47
37:DL:132:ARG:HA	37:DL:135:ILE:CG2	2.44	0.47
42:DN:101:GLY:O	42:DN:102:PHE:HB2	2.14	0.47
43:DO:66:GLY:O	43:DO:102:ARG:HD3	2.14	0.47
28:DP:16:VAL:HG23	28:DP:16:VAL:O	2.14	0.47
44:DQ:72:GLY:O	44:DQ:113:LYS:HE2	2.14	0.47
23:DB:572:A:H5''	49:DR:80:ARG:HH22	1.78	0.47
23:DB:483:A:H4'	46:DU:46:LYS:HA	1.97	0.47
39:DX:31:GLN:O	39:DX:37:LEU:HB2	2.13	0.47
1:AA:1465:A:H2'	1:AA:1466:C:C6	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:375:U:OP1	13:AP:70:ARG:HB2	2.13	0.47
1:AA:845:A:H5''	1:AA:846:G:H8	1.76	0.47
3:AD:25:ARG:O	3:AD:27:ILE:HG13	2.14	0.47
7:AH:29:SER:OG	7:AH:32:LYS:HG3	2.13	0.47
7:AH:17:GLN:OE1	7:AH:69:ALA:HB1	2.14	0.47
9:AJ:18:ILE:HG12	9:AJ:72:ARG:CG	2.44	0.47
10:AK:52:ARG:NH1	10:AK:52:ARG:HB3	2.29	0.47
21:AN:20:PHE:CD2	21:AN:55:SER:HA	2.49	0.47
14:AQ:24:ILE:HD13	14:AQ:43:LEU:HD13	1.95	0.47
19:AU:36:PHE:O	19:AU:39:LYS:HD2	2.14	0.47
31:B0:41:HIS:CB	42:BN:99:LYS:HB2	2.44	0.47
32:B4:36:ARG:HG2	32:B4:37:GLN:N	2.23	0.47
23:BB:1182:G:H2'	23:BB:1183:U:O4'	2.14	0.47
23:BB:1242:U:H2'	23:BB:1243:C:C6	2.49	0.47
23:BB:1387:A:H2'	23:BB:1388:G:C8	2.49	0.47
23:BB:1459:G:H2'	23:BB:1461:C:C4	2.49	0.47
23:BB:1563:U:O2'	23:BB:1564:C:H5'	2.14	0.47
23:BB:2135:A:H3'	23:BB:2136:G:H8	1.79	0.47
23:BB:227:A:C2	23:BB:2407:A:H1'	2.50	0.47
23:BB:2836:U:H2'	23:BB:2837:A:C8	2.50	0.47
23:BB:404:A:H4'	23:BB:405:U:O5'	2.15	0.47
23:BB:552:U:O2'	23:BB:553:G:H5'	2.15	0.47
23:BB:677:A:H2'	23:BB:678:C:H6	1.78	0.47
23:BB:955:U:H5''	38:BM:86:LYS:NZ	2.30	0.47
25:BC:202:ARG:HE	25:BC:213:ARG:HH21	1.61	0.47
25:BC:209:ALA:HA	25:BC:212:TRP:NE1	2.29	0.47
29:BE:16:GLU:O	29:BE:20:GLY:HA3	2.14	0.47
48:BG:129:GLU:C	48:BG:130:ILE:HG13	2.34	0.47
40:BH:73:ASN:N	40:BH:73:ASN:ND2	2.51	0.47
27:BK:35:VAL:CG2	27:BK:36:GLY:H	2.12	0.47
27:BK:51:LYS:O	27:BK:52:VAL:HG13	2.14	0.47
27:BK:58:LEU:HD11	27:BK:86:LEU:HB2	1.96	0.47
38:BM:73:ILE:HG21	38:BM:91:TYR:CZ	2.50	0.47
42:BN:90:ARG:HB3	42:BN:94:TYR:HE1	1.79	0.47
43:BO:115:LEU:H	43:BO:115:LEU:CD2	2.26	0.47
28:BP:3:ILE:HD13	28:BP:4:ILE:N	2.29	0.47
46:BU:28:LEU:HB2	46:BU:32:LYS:O	2.15	0.47
46:BU:73:ASN:C	46:BU:75:ALA:H	2.17	0.47
35:BV:42:LEU:HB2	35:BV:47:VAL:HG21	1.95	0.47
51:BZ:14:THR:HA	51:BZ:28:ARG:HA	1.96	0.47
1:CA:1053:G:C4'	1:CA:1054:C:H5'	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1155:A:O2'	1:CA:1156:G:H5'	2.14	0.47
1:CA:237:G:H5''	14:CQ:26:ARG:NH2	2.29	0.47
1:CA:451:A:H1'	1:CA:452:A:C8	2.49	0.47
1:CA:76:G:H2'	1:CA:77:A:H8	1.79	0.47
1:CA:94:G:H4'	1:CA:95:C:C5'	2.45	0.47
18:CB:20:ARG:HD2	18:CB:38:HIS:CE1	2.49	0.47
18:CB:213:LEU:HA	18:CB:213:LEU:HD23	1.79	0.47
18:CB:19:THR:O	18:CB:37:VAL:HA	2.14	0.47
4:CE:131:ASN:O	4:CE:135:VAL:HG23	2.14	0.47
8:CI:20:ILE:HD13	8:CI:85:ALA:CB	2.44	0.47
8:CI:56:MET:C	8:CI:58:GLU:N	2.67	0.47
10:CK:19:VAL:N	10:CK:34:THR:O	2.47	0.47
17:CT:70:LYS:HA	17:CT:73:ARG:CZ	2.45	0.47
19:CU:34:ARG:HD3	19:CU:39:LYS:CE	2.39	0.47
31:D0:46:GLY:O	31:D0:53:VAL:N	2.48	0.47
23:DB:1283:G:N2	23:DB:1285:A:H3'	2.29	0.47
23:DB:1439:A:H5''	23:DB:1440:U:OP2	2.14	0.47
23:DB:1869:G:N3	23:DB:1869:G:H2'	2.30	0.47
23:DB:1977:A:H2'	23:DB:1978:A:C8	2.49	0.47
23:DB:1998:A:H2'	23:DB:1999:C:H6	1.79	0.47
23:DB:2027:G:O2'	23:DB:2028:U:H5'	2.14	0.47
23:DB:2348:U:O2'	23:DB:2349:G:H5'	2.15	0.47
23:DB:2389:G:C5'	23:DB:2390:U:H5'	2.44	0.47
23:DB:407:G:O2'	23:DB:408:G:H5'	2.15	0.47
23:DB:919:U:H6	23:DB:919:U:O5'	1.97	0.47
26:DD:107:VAL:N	26:DD:206:ALA:H	2.12	0.47
47:DF:106:ALA:HA	47:DF:135:ILE:CD1	2.45	0.47
40:DH:26:ALA:O	40:DH:27:ARG:C	2.52	0.47
24:DI:100:ILE:O	24:DI:139:VAL:HG13	2.14	0.47
24:DI:35:MET:HE3	24:DI:39:LYS:HG2	1.96	0.47
43:DO:83:LEU:CD1	43:DO:114:GLY:HA3	2.44	0.47
28:DP:3:ILE:HG23	28:DP:4:ILE:N	2.30	0.47
44:DQ:59:LEU:O	44:DQ:62:ALA:HB3	2.14	0.47
23:DB:996:A:H4'	44:DQ:91:ARG:HG2	1.96	0.47
50:DT:87:LEU:HB2	50:DT:91:GLN:HG2	1.95	0.47
46:DU:41:VAL:O	46:DU:42:LYS:HB2	2.14	0.47
46:DU:46:LYS:HE3	46:DU:47:PRO:O	2.13	0.47
35:DV:29:ILE:HD13	35:DV:31:TYR:CE2	2.50	0.47
52:DW:13:ARG:HH11	52:DW:13:ARG:HA	1.79	0.47
51:DZ:40:VAL:O	51:DZ:42:SER:N	2.46	0.47
1:AA:1336:C:H4'	1:AA:1337:G:N3	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1532:U:H2'	1:AA:1533:C:H3'	1.95	0.47
18:AB:61:SER:HA	18:AB:224:ARG:HA	1.97	0.47
18:AB:68:PHE:HA	18:AB:161:PHE:O	2.13	0.47
3:AD:24:VAL:O	3:AD:27:ILE:HD11	2.14	0.47
4:AE:17:VAL:HG23	4:AE:33:THR:O	2.14	0.47
4:AE:36:THR:HB	4:AE:63:MET:HE2	1.95	0.47
8:AI:64:ILE:HD12	8:AI:64:ILE:N	2.30	0.47
10:AK:14:GLN:OE1	10:AK:77:GLY:HA3	2.14	0.47
12:AM:78:ARG:HH11	12:AM:78:ARG:HG2	1.79	0.47
1:AA:751:U:H4'	20:AO:24:SER:HA	1.96	0.47
16:AS:9:PHE:HE1	16:AS:36:ARG:HG3	1.79	0.47
33:B1:50:GLU:O	33:B1:51:ALA:HB2	2.15	0.47
22:BA:29:A:OP1	43:BO:31:THR:HB	2.14	0.47
23:BB:1026:G:OP2	23:BB:1134:A:H1'	2.13	0.47
23:BB:1158:C:H5'	30:BY:31:ILE:HG12	1.96	0.47
23:BB:1290:C:O2'	23:BB:1291:C:H5'	2.14	0.47
23:BB:136:G:H2'	23:BB:137:U:O4'	2.14	0.47
23:BB:138:U:O2'	23:BB:141:G:N2	2.47	0.47
23:BB:1461:C:O2'	23:BB:1462:C:H5'	2.13	0.47
23:BB:1921:G:O2'	23:BB:1922:G:H5'	2.14	0.47
23:BB:2029:G:H2'	23:BB:2031:A:OP1	2.15	0.47
23:BB:2134:A:H2'	23:BB:2135:A:C8	2.50	0.47
23:BB:2252:G:O2'	23:BB:2253:G:H5'	2.13	0.47
23:BB:2504:U:O5'	23:BB:2504:U:H6	1.97	0.47
23:BB:2654:A:N1	23:BB:2665:A:H5''	2.30	0.47
23:BB:2807:U:H5'	23:BB:2808:G:OP2	2.14	0.47
23:BB:457:A:N1	23:BB:470:A:H5''	2.30	0.47
23:BB:27:G:H1'	23:BB:513:A:N6	2.29	0.47
23:BB:567:U:H2'	23:BB:568:U:O4'	2.14	0.47
23:BB:634:C:O2'	23:BB:635:C:H5'	2.14	0.47
23:BB:945:A:H4'	23:BB:945:A:OP2	2.14	0.47
25:BC:173:LEU:HD22	25:BC:181:ARG:O	2.14	0.47
26:BD:106:LYS:HB3	26:BD:206:ALA:HB3	1.95	0.47
26:BD:8:LYS:HG3	26:BD:9:VAL:N	2.30	0.47
47:BF:163:GLU:O	47:BF:166:ARG:HB2	2.14	0.47
48:BG:26:LYS:HA	48:BG:32:LEU:HA	1.96	0.47
48:BG:71:LEU:HA	48:BG:74:MET:SD	2.54	0.47
40:BH:103:VAL:HB	40:BH:108:VAL:HB	1.96	0.47
40:BH:81:ALA:CB	40:BH:146:VAL:HA	2.44	0.47
27:BK:115:ILE:HG23	27:BK:116:ILE:H	1.80	0.47
43:BO:31:THR:HG23	43:BO:34:HIS:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:15:PHE:N	28:BP:11:GLN:HE22	2.12	0.47
46:BU:47:PRO:HD3	46:BU:55:GLY:CA	2.45	0.47
35:BV:29:ILE:HG13	35:BV:88:HIS:CE1	2.49	0.47
35:BV:42:LEU:HD12	35:BV:47:VAL:HG21	1.96	0.47
35:BV:68:LYS:HD3	35:BV:68:LYS:N	2.28	0.47
1:CA:18:C:C1'	1:CA:1079:G:H21	2.27	0.47
1:CA:1495:U:H2'	1:CA:1496:C:C6	2.49	0.47
1:CA:599:C:O2'	1:CA:600:A:H5'	2.14	0.47
18:CB:107:ARG:HA	18:CB:110:ILE:CG1	2.44	0.47
18:CB:164:ASP:HB2	18:CB:203:ASP:HB2	1.97	0.47
2:CC:104:GLU:HG2	2:CC:105:VAL:H	1.80	0.47
7:CH:124:ILE:HG22	7:CH:125:ILE:H	1.78	0.47
8:CI:42:THR:O	8:CI:46:VAL:HG13	2.13	0.47
21:CN:77:GLY:O	21:CN:78:LEU:HD12	2.15	0.47
20:CO:33:THR:HG23	20:CO:63:ARG:HH11	1.79	0.47
14:CQ:80:LYS:HD2	14:CQ:80:LYS:C	2.35	0.47
31:D0:12:ARG:HD2	31:D0:16:ARG:NH1	2.30	0.47
31:D0:27:LEU:HD22	31:D0:38:LEU:HA	1.95	0.47
23:DB:1592:C:H2'	23:DB:1593:A:H8	1.80	0.47
23:DB:2203:U:C2'	23:DB:2204:G:OP2	2.63	0.47
23:DB:2324:U:H5'	23:DB:2325:G:H5''	1.95	0.47
23:DB:1664:A:H1'	23:DB:2726:A:C2	2.49	0.47
23:DB:754:U:H2'	23:DB:755:U:C6	2.50	0.47
23:DB:765:C:O2'	23:DB:766:U:H5'	2.14	0.47
23:DB:875:G:H2'	23:DB:876:C:O4'	2.14	0.47
23:DB:973:A:OP1	23:DB:973:A:H8	1.97	0.47
23:DB:1820:U:N3	25:DC:197:ALA:O	2.47	0.47
25:DC:255:LYS:C	25:DC:256:THR:HG23	2.33	0.47
25:DC:2:VAL:HG23	25:DC:3:VAL:N	2.22	0.47
26:DD:7:LYS:HG3	26:DD:198:GLY:O	2.13	0.47
47:DF:105:ILE:HA	47:DF:108:PRO:HB2	1.95	0.47
48:DG:112:VAL:O	48:DG:113:ASP:HB2	2.14	0.47
48:DG:166:GLU:N	48:DG:166:GLU:CD	2.65	0.47
24:DI:18:ASN:HB2	24:DI:38:CYS:SG	2.54	0.47
41:DJ:56:VAL:HG12	41:DJ:57:LEU:N	2.29	0.47
41:DJ:6:ALA:HB3	41:DJ:45:THR:CG2	2.42	0.47
37:DL:78:ARG:CB	37:DL:113:ALA:HB2	2.38	0.47
37:DL:74:THR:HA	37:DL:107:PHE:O	2.15	0.47
38:DM:34:LYS:O	38:DM:128:THR:HG22	2.14	0.47
38:DM:38:ARG:HG2	38:DM:98:PRO:HD3	1.96	0.47
28:DP:3:ILE:O	28:DP:7:LEU:HD13	2.12	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DS:31:GLN:O	45:DS:34:ASP:HB2	2.13	0.47
46:DU:78:LYS:HE3	46:DU:79:ALA:N	2.30	0.47
39:DX:20:ASN:N	39:DX:20:ASN:HD22	2.09	0.47
30:DY:2:LYS:HE2	30:DY:39:ASP:OD2	2.14	0.47
1:AA:1293:C:H2'	1:AA:1294:G:C8	2.50	0.47
1:AA:438:U:H4'	3:AD:119:HIS:HD2	1.78	0.47
1:AA:525:C:H2'	1:AA:526:C:C6	2.49	0.47
1:AA:948:C:C2'	1:AA:949:A:H5'	2.45	0.47
18:AB:77:GLU:HA	18:AB:80:LYS:HD3	1.96	0.47
2:AC:112:ALA:O	2:AC:113:LYS:C	2.52	0.47
2:AC:129:PHE:CB	2:AC:133:MET:HE3	2.44	0.47
2:AC:45:GLU:C	2:AC:46:LEU:HD22	2.35	0.47
2:AC:52:SER:OG	2:AC:53:ARG:N	2.47	0.47
1:AA:405:U:C5	3:AD:4:LEU:HD21	2.48	0.47
9:AJ:33:GLY:O	9:AJ:34:ALA:C	2.53	0.47
1:AA:950:U:H3'	12:AM:100:ARG:NH2	2.30	0.47
20:AO:39:LEU:CD2	20:AO:56:LEU:HB2	2.44	0.47
16:AS:29:PRO:HB3	16:AS:47:THR:C	2.35	0.47
16:AS:39:ILE:HD12	16:AS:65:MET:O	2.15	0.47
17:AT:24:ARG:CG	17:AT:65:LEU:HD11	2.45	0.47
22:BA:75:G:H1	22:BA:102:G:N2	2.12	0.47
23:BB:1292:G:O2'	23:BB:1293:C:H5'	2.15	0.47
23:BB:1459:G:H3'	23:BB:1460:U:C4'	2.44	0.47
23:BB:1576:U:O2'	23:BB:1577:C:H5'	2.14	0.47
23:BB:1625:C:H3'	23:BB:1626:A:C8	2.50	0.47
23:BB:1744:A:H3'	23:BB:1745:A:C8	2.49	0.47
23:BB:1917:U:H2'	23:BB:1918:A:H8	1.79	0.47
23:BB:1131:G:N2	23:BB:2024:G:H21	2.13	0.47
23:BB:2210:U:N3	23:BB:2212:A:N7	2.62	0.47
23:BB:234:U:O2'	23:BB:235:U:H5'	2.14	0.47
23:BB:2365:G:O2'	52:BW:59:PHE:CE1	2.68	0.47
23:BB:2454:G:C2'	23:BB:2455:G:H5'	2.44	0.47
23:BB:2634:A:H2'	23:BB:2635:A:C8	2.49	0.47
23:BB:2809:A:H2'	23:BB:2810:A:C8	2.49	0.47
23:BB:2821:A:H3'	23:BB:2821:A:OP2	2.14	0.47
23:BB:400:G:N7	51:BZ:57:ARG:NH1	2.62	0.47
23:BB:68:G:H2'	23:BB:69:C:H6	1.79	0.47
23:BB:796:C:H2'	23:BB:797:G:C8	2.49	0.47
23:BB:802:A:H2'	23:BB:803:U:H6	1.79	0.47
23:BB:863:A:H2'	23:BB:864:G:C8	2.48	0.47
25:BC:250:GLN:C	25:BC:252:LYS:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2580:U:H5'	26:BD:136:ASN:H	1.79	0.47
26:BD:54:ALA:N	26:BD:76:GLY:HA2	2.30	0.47
23:BB:659:G:H21	29:BE:30:GLN:HE22	1.61	0.47
48:BG:152:ARG:HD3	48:BG:153:PRO:CD	2.42	0.47
24:BI:32:VAL:HG13	24:BI:66:PHE:CD2	2.49	0.47
37:BL:19:LEU:HD22	37:BL:31:GLY:O	2.13	0.47
1:AA:1441:A:H2	28:BP:113:LEU:HD22	1.79	0.47
44:BQ:71:ASN:HD22	44:BQ:109:VAL:HG21	1.79	0.47
35:BV:51:GLN:NE2	35:BV:79:ARG:HH22	2.13	0.47
1:CA:1234:C:H1'	1:CA:1364:U:O2	2.14	0.47
1:CA:435:A:H2'	1:CA:435:A:N3	2.29	0.47
1:CA:552:U:H4'	11:CL:82:ARG:HG2	1.97	0.47
1:CA:980:C:H4'	21:CN:12:ARG:HH22	1.80	0.47
18:CB:118:THR:O	18:CB:119:GLN:C	2.51	0.47
3:CD:146:GLU:HA	3:CD:149:LYS:CG	2.41	0.47
3:CD:170:LEU:HD12	3:CD:170:LEU:O	2.15	0.47
4:CE:82:HIS:HB2	4:CE:83:PRO:HD2	1.96	0.47
6:CG:106:ALA:O	6:CG:110:ARG:HG3	2.14	0.47
8:CI:108:ARG:HB3	8:CI:108:ARG:HH11	1.79	0.47
8:CI:74:GLN:O	8:CI:78:ILE:HG13	2.15	0.47
9:CJ:28:THR:O	9:CJ:31:ARG:HG2	2.13	0.47
14:CQ:37:ILE:HG22	14:CQ:38:LYS:N	2.29	0.47
14:CQ:10:ARG:HA	14:CQ:56:ASP:O	2.14	0.47
14:CQ:80:LYS:H	14:CQ:80:LYS:NZ	2.13	0.47
17:CT:72:ALA:HA	17:CT:75:LYS:HD3	1.96	0.47
34:D3:21:PHE:HB2	34:D3:49:VAL:HG22	1.97	0.47
23:DB:103:A:H3'	23:DB:104:A:C8	2.48	0.47
23:DB:115:C:O2'	23:DB:116:C:H5'	2.14	0.47
23:DB:1530:G:H2'	23:DB:1530:G:N3	2.29	0.47
23:DB:1563:U:O2'	23:DB:1564:C:H5'	2.15	0.47
23:DB:1690:A:H2'	23:DB:1691:C:O4'	2.15	0.47
23:DB:1864:U:O2'	23:DB:1865:U:H5'	2.14	0.47
23:DB:2455:G:H2'	23:DB:2456:C:H6	1.79	0.47
23:DB:2559:C:H2'	23:DB:2560:A:H8	1.80	0.47
23:DB:264:C:H2'	23:DB:265:A:C5'	2.43	0.47
23:DB:418:C:H2'	23:DB:419:U:C6	2.49	0.47
23:DB:423:A:H5''	23:DB:424:G:C5'	2.44	0.47
23:DB:425:G:O2'	23:DB:426:C:H5'	2.14	0.47
23:DB:557:C:H2'	23:DB:558:U:C6	2.50	0.47
23:DB:596:U:O2'	23:DB:597:G:H5'	2.14	0.47
23:DB:784:G:H5''	25:DC:225:ASN:OD1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:855:G:H21	52:DW:23:LYS:CG	2.18	0.47
26:DD:46:ARG:HG3	26:DD:46:ARG:NH1	2.30	0.47
22:DA:54:G:N2	47:DF:25:MET:HG2	2.28	0.47
40:DH:110:VAL:HG23	40:DH:132:PHE:CE2	2.49	0.47
24:DI:11:GLN:HA	24:DI:55:PRO:HA	1.95	0.47
24:DI:129:GLU:CB	24:DI:133:ARG:HH12	2.27	0.47
24:DI:49:GLU:HB3	24:DI:52:LEU:HD12	1.97	0.47
24:DI:72:THR:CG2	24:DI:112:LYS:HD2	2.44	0.47
41:DJ:13:ARG:HB3	41:DJ:53:TYR:CD2	2.49	0.47
41:DJ:34:ARG:HG3	41:DJ:39:LYS:HB3	1.96	0.47
37:DL:76:GLU:HG3	37:DL:111:ILE:HG13	1.96	0.47
37:DL:78:ARG:HB3	37:DL:113:ALA:CB	2.37	0.47
42:DN:19:ALA:C	42:DN:21:PHE:N	2.68	0.47
42:DN:37:THR:CG2	42:DN:39:PRO:HD2	2.34	0.47
49:DR:38:VAL:HG22	49:DR:40:MET:H	1.78	0.47
46:DU:47:PRO:HD3	46:DU:55:GLY:CA	2.45	0.47
30:DY:47:ILE:HG21	30:DY:56:VAL:CG2	2.45	0.47
51:DZ:21:ALA:HB3	51:DZ:23:ASN:HD21	1.79	0.47
51:DZ:3:ARG:HG3	51:DZ:50:ARG:HH11	1.78	0.47
1:AA:1027:C:H2'	1:AA:1028:C:C6	2.50	0.47
1:AA:1127:G:H5'	1:AA:1280:A:O2'	2.14	0.47
1:AA:410:G:P	3:AD:25:ARG:HD2	2.54	0.47
1:AA:547:A:H4'	1:AA:548:G:O5'	2.15	0.47
1:AA:635:A:H2'	1:AA:636:U:C6	2.50	0.47
1:AA:818:G:C3'	1:AA:819:A:H5''	2.43	0.47
1:AA:842:U:H6	1:AA:842:U:OP1	1.97	0.47
18:AB:145:ASN:ND2	18:AB:145:ASN:N	2.62	0.47
3:AD:137:SER:HB2	3:AD:138:PRO:HD2	1.96	0.47
5:AF:93:LYS:O	5:AF:94:HIS:HB2	2.14	0.47
7:AH:55:LYS:CE	7:AH:55:LYS:HA	2.43	0.47
1:AA:1152:A:OP1	9:AJ:70:HIS:ND1	2.47	0.47
12:AM:93:GLY:HA2	12:AM:108:ARG:HH12	1.80	0.47
14:AQ:7:LEU:O	14:AQ:60:ILE:HD13	2.14	0.47
16:AS:64:GLU:N	16:AS:64:GLU:CD	2.68	0.47
31:B0:2:VAL:HG12	31:B0:3:GLN:N	2.30	0.47
23:BB:1197:G:O2'	23:BB:1198:U:H5'	2.15	0.47
23:BB:1390:U:O2'	23:BB:1391:U:H5'	2.15	0.47
23:BB:1523:U:H5''	23:BB:1524:G:C8	2.49	0.47
23:BB:151:C:H2'	23:BB:152:A:H8	1.79	0.47
23:BB:1542:U:H2'	23:BB:1543:G:O4'	2.14	0.47
23:BB:1582:C:H2'	23:BB:1583:A:O4'	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:15:G:O2'	23:BB:16:C:H5'	2.14	0.47
23:BB:1859:U:H2'	23:BB:1860:G:H8	1.77	0.47
23:BB:189:G:H2'	23:BB:205:G:N2	2.30	0.47
23:BB:230:G:H2'	23:BB:231:A:H8	1.79	0.47
23:BB:2362:C:O2'	23:BB:2363:G:H5'	2.14	0.47
23:BB:2722:G:H2'	23:BB:2723:C:O4'	2.14	0.47
23:BB:638:G:H2'	23:BB:639:U:C6	2.50	0.47
23:BB:691:C:H5'	25:BC:216:ARG:NH1	2.30	0.47
23:BB:779:U:O2'	23:BB:780:G:H5'	2.15	0.47
25:BC:169:ALA:O	25:BC:185:ALA:HB3	2.15	0.47
25:BC:239:PHE:C	25:BC:241:LYS:H	2.17	0.47
25:BC:255:LYS:C	25:BC:256:THR:HG23	2.35	0.47
25:BC:93:VAL:CG1	25:BC:101:ARG:H	2.28	0.47
26:BD:109:VAL:HG11	26:BD:193:VAL:HB	1.96	0.47
26:BD:154:LYS:HG2	26:BD:155:VAL:N	2.30	0.47
29:BE:129:PRO:HB3	29:BE:159:LEU:HD23	1.97	0.47
41:BJ:38:GLY:O	41:BJ:43:GLU:HB2	2.14	0.47
27:BK:60:ALA:HA	27:BK:87:LEU:HD22	1.95	0.47
37:BL:95:LEU:O	37:BL:100:ILE:HG22	2.14	0.47
37:BL:18:ARG:O	37:BL:19:LEU:HD12	2.14	0.47
43:BO:17:LYS:O	43:BO:21:LEU:HB2	2.14	0.47
41:BJ:44:TYR:CE2	44:BQ:59:LEU:HD11	2.50	0.47
35:BV:59:GLU:HG2	35:BV:60:VAL:H	1.79	0.47
1:CA:1061:G:H2'	1:CA:1062:U:H6	1.80	0.47
1:CA:1110:A:H2'	1:CA:1111:A:H8	1.80	0.47
1:CA:518:C:H2'	1:CA:530:G:N7	2.30	0.47
1:CA:640:A:O2'	1:CA:641:U:H5'	2.15	0.47
1:CA:994:A:C2	21:CN:4:SER:HA	2.49	0.47
1:CA:996:A:H2'	1:CA:997:U:C6	2.49	0.47
18:CB:214:GLY:C	18:CB:216:VAL:H	2.17	0.47
18:CB:46:VAL:HA	18:CB:49:PHE:HD2	1.78	0.47
3:CD:197:HIS:O	3:CD:201:GLU:HG3	2.15	0.47
3:CD:90:LEU:HD21	3:CD:196:GLU:CB	2.45	0.47
4:CE:132:PRO:HG2	4:CE:133:ILE:H	1.79	0.47
1:CA:778:G:H21	10:CK:121:ARG:HD3	1.79	0.47
1:CA:255:G:H1'	14:CQ:17:GLU:OE2	2.14	0.47
15:CR:52:ARG:O	15:CR:56:ARG:HG3	2.14	0.47
31:D0:18:HIS:C	31:D0:20:ALA:H	2.17	0.47
23:DB:1092:C:C2'	23:DB:1093:G:H5'	2.44	0.47
23:DB:1150:C:H2'	23:DB:1151:A:C8	2.49	0.47
23:DB:1324:G:H3'	23:DB:1325:U:H4'	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1387:A:H2'	23:DB:1388:G:C8	2.49	0.47
23:DB:1808:A:H3'	23:DB:1809:A:H8	1.79	0.47
23:DB:2053:G:O2'	23:DB:2054:A:H5'	2.14	0.47
23:DB:21:A:H2'	23:DB:22:C:H6	1.78	0.47
23:DB:2260:C:O2'	23:DB:2261:C:H5'	2.14	0.47
23:DB:2294:G:P	43:DO:94:ARG:HH12	2.37	0.47
23:DB:233:A:H61	23:DB:428:A:H61	1.62	0.47
23:DB:2471:A:O2'	23:DB:2472:G:O5'	2.33	0.47
23:DB:277:G:H2'	23:DB:278:A:OP2	2.14	0.47
23:DB:346:A:H5'	23:DB:346:A:N3	2.29	0.47
23:DB:543:G:C2'	23:DB:545:U:H5'	2.44	0.47
23:DB:558:U:H5'	41:DJ:114:LEU:HD22	1.96	0.47
25:DC:115:ILE:HB	25:DC:126:GLY:O	2.14	0.47
25:DC:270:ARG:HB3	25:DC:270:ARG:NH1	2.29	0.47
26:DD:204:LYS:NZ	26:DD:204:LYS:HB3	2.28	0.47
26:DD:29:VAL:O	26:DD:185:ASN:HB3	2.15	0.47
26:DD:5:VAL:H	26:DD:32:ASN:HD21	1.61	0.47
47:DF:40:GLY:HA2	47:DF:84:ILE:O	2.14	0.47
48:DG:7:PRO:O	48:DG:8:VAL:HB	2.14	0.47
40:DH:121:VAL:HG23	40:DH:122:LEU:N	2.29	0.47
41:DJ:124:VAL:O	41:DJ:125:TYR:HB2	2.15	0.47
27:DK:109:SER:O	27:DK:111:LYS:N	2.47	0.47
27:DK:88:ASN:HD22	27:DK:89:ASN:H	1.62	0.47
38:DM:83:GLY:O	38:DM:84:LYS:HG2	2.14	0.47
42:DN:101:GLY:CA	42:DN:109:PRO:HA	2.44	0.47
42:DN:2:ARG:HG2	42:DN:5:LYS:CB	2.38	0.47
42:DN:49:GLU:HA	42:DN:94:TYR:HD2	1.80	0.47
28:DP:4:ILE:C	28:DP:6:GLN:H	2.18	0.47
49:DR:41:ILE:HD13	49:DR:41:ILE:N	2.29	0.47
45:DS:4:ILE:HG21	45:DS:106:VAL:HG22	1.96	0.47
45:DS:30:SER:HA	45:DS:33:LEU:HD12	1.96	0.47
1:AA:1152:A:H2'	1:AA:1153:G:H8	1.80	0.47
1:AA:244:U:O4	1:AA:906:A:H1'	2.14	0.47
1:AA:202:G:N2	1:AA:465:A:H61	2.10	0.47
18:AB:17:HIS:CG	18:AB:18:GLN:N	2.83	0.47
3:AD:41:GLY:C	3:AD:43:ARG:H	2.18	0.47
4:AE:104:ILE:HD11	4:AE:111:ARG:HA	1.96	0.47
5:AF:11:HIS:HB3	5:AF:14:GLN:CG	2.38	0.47
7:AH:14:ARG:HE	7:AH:75:GLN:NE2	2.13	0.47
7:AH:82:LEU:O	7:AH:82:LEU:HD13	2.15	0.47
10:AK:21:HIS:HA	10:AK:84:MET:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:80:LEU:HB3	11:AL:97:VAL:CG2	2.45	0.47
11:AL:85:ARG:HA	11:AL:93:ARG:HA	1.96	0.47
20:AO:36:ILE:HD12	20:AO:60:VAL:CG2	2.45	0.47
14:AQ:24:ILE:HD12	14:AQ:24:ILE:N	2.30	0.47
15:AR:20:ILE:HG22	15:AR:53:GLN:CD	2.34	0.47
16:AS:43:MET:HE3	16:AS:61:VAL:HG21	1.97	0.47
1:AA:263:A:OP1	17:AT:73:ARG:NH1	2.48	0.47
23:BB:2284:A:OP1	33:B1:3:GLY:O	2.33	0.47
23:BB:1266:G:N2	23:BB:2012:G:H2'	2.30	0.47
23:BB:1636:U:H2'	23:BB:1637:A:H8	1.78	0.47
23:BB:1818:U:HO2'	23:BB:1819:A:P	2.37	0.47
23:BB:213:A:O2'	23:BB:214:G:H5'	2.15	0.47
23:BB:241:A:OP1	23:BB:241:A:H8	1.97	0.47
23:BB:2895:G:H2'	23:BB:2896:C:C6	2.50	0.47
23:BB:718:A:H5'	23:BB:719:C:OP2	2.15	0.47
23:BB:998:C:H3'	56:BB:3247:HOH:O	2.15	0.47
23:BB:1567:G:H2'	25:BC:82:TYR:HE1	1.79	0.47
29:BE:111:GLU:HA	29:BE:114:ARG:CG	2.35	0.47
29:BE:158:PHE:HA	29:BE:169:VAL:CG2	2.41	0.47
47:BF:41:GLU:CB	47:BF:48:LEU:HD11	2.44	0.47
48:BG:84:LYS:HB2	48:BG:132:LEU:N	2.30	0.47
48:BG:84:LYS:HG3	48:BG:131:VAL:CB	2.44	0.47
40:BH:18:GLN:HE21	40:BH:39:ALA:HB1	1.79	0.47
24:BI:122:GLU:CD	24:BI:122:GLU:H	2.17	0.47
41:BJ:65:THR:HG23	41:BJ:66:GLY:N	2.29	0.47
41:BJ:84:ILE:HG23	41:BJ:84:ILE:O	2.14	0.47
27:BK:47:ILE:CG1	27:BK:48:PRO:HD2	2.38	0.47
44:BQ:40:LYS:O	44:BQ:43:GLN:HB2	2.15	0.47
49:BR:28:ALA:HB3	49:BR:31:GLU:HG3	1.95	0.47
45:BS:32:ALA:O	45:BS:35:ILE:HG12	2.15	0.47
45:BS:35:ILE:HG22	45:BS:39:THR:CG2	2.44	0.47
23:BB:141:G:C6	50:BT:1:MET:HA	2.50	0.47
50:BT:62:VAL:HB	50:BT:64:LYS:HZ1	1.79	0.47
46:BU:48:VAL:O	46:BU:48:VAL:HG13	2.15	0.47
52:BW:43:LYS:O	52:BW:58:LEU:HD11	2.15	0.47
30:BY:6:ILE:HG22	30:BY:56:VAL:HA	1.97	0.47
1:CA:1137:C:O2	1:CA:1138:G:N1	2.47	0.47
1:CA:1053:G:N7	1:CA:1199:U:H2'	2.30	0.47
1:CA:960:U:O2'	1:CA:1223:C:H4'	2.14	0.47
1:CA:1296:C:H4'	1:CA:1302:C:N4	2.27	0.47
1:CA:1329:A:O2'	1:CA:1330:U:H5'	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1509:C:H2'	1:CA:1510:C:H6	1.80	0.47
1:CA:373:A:H3'	1:CA:373:A:P	2.55	0.47
1:CA:412:A:H4'	1:CA:413:G:OP1	2.15	0.47
1:CA:425:G:O2'	1:CA:426:U:H5'	2.15	0.47
1:CA:488:C:O2'	1:CA:489:C:H5'	2.15	0.47
1:CA:1057:G:H5''	2:CC:153:SER:OG	2.14	0.47
2:CC:70:ALA:HB2	2:CC:114:LEU:HD13	1.95	0.47
3:CD:137:SER:HB2	3:CD:138:PRO:HD2	1.96	0.47
3:CD:24:VAL:O	3:CD:27:ILE:HD11	2.15	0.47
6:CG:112:ASP:CB	6:CG:118:ARG:HG2	2.45	0.47
6:CG:91:ARG:HB3	6:CG:92:PRO:HD2	1.97	0.47
12:CM:100:ARG:O	12:CM:100:ARG:HG3	2.14	0.47
12:CM:38:ILE:HG22	12:CM:42:VAL:HG11	1.96	0.47
33:D1:8:ILE:HD12	33:D1:9:LYS:H	1.80	0.47
23:DB:1234:U:H2'	23:DB:1235:G:O4'	2.15	0.47
23:DB:1238:G:O2'	23:DB:1239:G:H5'	2.14	0.47
23:DB:1361:G:C6	23:DB:1371:G:N2	2.83	0.47
23:DB:1432:G:O2'	23:DB:1433:A:H5'	2.15	0.47
23:DB:1459:G:H3'	23:DB:1460:U:C4'	2.45	0.47
23:DB:1576:U:O2'	23:DB:1577:C:H5'	2.14	0.47
23:DB:1744:A:H3'	23:DB:1745:A:C8	2.49	0.47
23:DB:1680:U:O2	23:DB:1763:G:H3'	2.15	0.47
23:DB:1779:U:C5	23:DB:1784:A:N7	2.83	0.47
23:DB:1793:C:H2'	23:DB:1794:A:C8	2.49	0.47
23:DB:185:G:H2'	23:DB:186:G:O4'	2.14	0.47
23:DB:1945:G:H2'	23:DB:1946:U:C6	2.49	0.47
23:DB:19:A:H2'	23:DB:20:C:H6	1.79	0.47
23:DB:2504:U:H6	23:DB:2504:U:O5'	1.98	0.47
23:DB:2580:U:H5'	26:DD:136:ASN:H	1.80	0.47
23:DB:634:C:O2'	23:DB:635:C:H5'	2.15	0.47
23:DB:738:G:H2'	23:DB:739:A:C8	2.49	0.47
23:DB:852:U:H2'	23:DB:853:C:C6	2.50	0.47
23:DB:876:C:H42	23:DB:901:C:H42	1.59	0.47
26:DD:97:SER:OG	26:DD:98:VAL:N	2.48	0.47
47:DF:57:ALA:HB2	47:DF:64:PRO:CG	2.45	0.47
48:DG:154:GLU:C	48:DG:156:TYR:H	2.18	0.47
48:DG:26:LYS:HA	48:DG:32:LEU:HA	1.96	0.47
48:DG:51:PHE:CD2	48:DG:68:ARG:HG2	2.50	0.47
37:DL:61:LEU:N	37:DL:61:LEU:CD1	2.78	0.47
42:DN:28:LEU:N	42:DN:34:ILE:HD11	2.29	0.47
42:DN:37:THR:HG22	42:DN:39:PRO:CD	2.33	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DU:95:PHE:HB2	46:DU:100:GLU:HB3	1.97	0.47
35:DV:59:GLU:HG2	35:DV:60:VAL:H	1.78	0.47
51:DZ:32:ASN:O	51:DZ:33:LEU:C	2.53	0.47
1:AA:1041:G:O5'	1:AA:1041:G:H8	1.97	0.47
1:AA:1237:C:H5''	1:AA:1238:A:O4'	2.14	0.47
1:AA:1262:C:H42	1:AA:1273:C:H42	1.63	0.47
1:AA:1497:G:H2'	1:AA:1498:U:H5'	1.96	0.47
1:AA:211:G:H3'	1:AA:211:G:N3	2.30	0.47
1:AA:607:A:H2'	1:AA:608:A:C8	2.50	0.47
1:AA:889:A:H5'	1:AA:891:U:H1'	1.97	0.47
18:AB:24:PRO:C	18:AB:26:MET:H	2.18	0.47
2:AC:55:VAL:HG12	2:AC:56:ILE:N	2.30	0.47
5:AF:9:MET:HA	5:AF:58:HIS:O	2.14	0.47
7:AH:124:ILE:HG22	7:AH:125:ILE:H	1.78	0.47
12:AM:15:VAL:HG22	12:AM:29:SER:HB3	1.97	0.47
23:BB:1324:G:H3'	23:BB:1325:U:H4'	1.96	0.47
23:BB:1528:A:H2'	23:BB:1529:G:O4'	2.15	0.47
23:BB:2010:G:H2'	23:BB:2011:U:H6	1.79	0.47
23:BB:2088:A:H2'	23:BB:2089:C:C6	2.50	0.47
23:BB:2634:A:H2'	23:BB:2635:A:H8	1.79	0.47
23:BB:2699:C:H2'	23:BB:2700:A:H8	1.78	0.47
23:BB:2757:A:H2	48:BG:63:GLN:NE2	2.12	0.47
23:BB:2813:A:H2'	23:BB:2814:A:H8	1.80	0.47
23:BB:909:A:H2'	23:BB:912:C:C5	2.49	0.47
25:BC:103:ILE:HG22	25:BC:104:LEU:N	2.30	0.47
23:BB:2580:U:C5'	26:BD:136:ASN:H	2.27	0.47
26:BD:202:ILE:HD12	26:BD:202:ILE:N	2.30	0.47
29:BE:67:ARG:HG2	29:BE:67:ARG:NH1	2.28	0.47
47:BF:167:ALA:O	47:BF:170:ALA:HB3	2.15	0.47
47:BF:31:GLU:HB3	47:BF:157:THR:HA	1.96	0.47
23:BB:2314:A:H4'	47:BF:34:THR:HG21	1.96	0.47
40:BH:116:ARG:HB2	40:BH:131:SER:HB2	1.96	0.47
27:BK:109:SER:O	27:BK:111:LYS:N	2.48	0.47
42:BN:35:LYS:HA	42:BN:111:ALA:O	2.15	0.47
42:BN:55:ALA:HB2	42:BN:79:LEU:HD22	1.96	0.47
50:BT:39:THR:HG22	50:BT:42:GLU:CG	2.44	0.47
46:BU:3:LYS:HD3	46:BU:82:VAL:HG21	1.96	0.47
1:CA:1270:G:H2'	1:CA:1271:A:C8	2.49	0.47
1:CA:191:G:H2'	1:CA:192:A:C8	2.49	0.47
1:CA:382:A:H2'	1:CA:383:A:O4'	2.15	0.47
1:CA:420:U:O2'	1:CA:421:U:H5''	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:68:G:C2	1:CA:69:G:H1'	2.50	0.47
1:CA:780:A:O2'	1:CA:781:A:H5''	2.15	0.47
18:CB:40:ILE:HG22	18:CB:41:ASN:N	2.30	0.47
18:CB:50:ASN:O	18:CB:51:GLU:C	2.53	0.47
4:CE:103:GLY:O	4:CE:121:ASN:HA	2.15	0.47
6:CG:46:LEU:HG	6:CG:57:GLU:HG2	1.97	0.47
9:CJ:98:VAL:C	9:CJ:100:ILE:H	2.18	0.47
10:CK:41:LEU:HD22	10:CK:76:TYR:CD2	2.49	0.47
12:CM:68:LEU:O	12:CM:72:ILE:HG22	2.15	0.47
14:CQ:7:LEU:HD22	14:CQ:24:ILE:HG21	1.96	0.47
14:CQ:46:HIS:HB2	14:CQ:66:LEU:HD13	1.96	0.47
23:DB:1128:G:O2'	23:DB:1129:A:H5''	2.14	0.47
23:DB:1182:G:H2'	23:DB:1183:U:O4'	2.14	0.47
23:DB:1665:A:H2'	23:DB:1666:G:O4'	2.14	0.47
23:DB:1672:A:C6	23:DB:1673:G:C6	3.02	0.47
23:DB:1839:G:C8	23:DB:1927:A:H1'	2.50	0.47
23:DB:2254:C:H2'	23:DB:2255:G:O4'	2.13	0.47
23:DB:2363:G:O2'	23:DB:2364:C:H5'	2.14	0.47
23:DB:299:A:H2'	23:DB:300:A:C8	2.48	0.47
23:DB:962:G:H2'	23:DB:963:U:C6	2.50	0.47
26:DD:121:THR:HG21	26:DD:127:PHE:CE1	2.50	0.47
26:DD:1:MET:SD	26:DD:2:ILE:N	2.85	0.47
47:DF:117:SER:HB3	47:DF:120:SER:HB3	1.95	0.47
47:DF:134:GLN:HB3	47:DF:149:ARG:HB3	1.97	0.47
47:DF:163:GLU:O	47:DF:166:ARG:HB2	2.14	0.47
24:DI:18:ASN:N	24:DI:19:PRO:CD	2.77	0.47
42:DN:51:LEU:O	42:DN:54:LEU:HB3	2.14	0.47
42:DN:55:ALA:HB2	42:DN:79:LEU:HD22	1.96	0.47
28:DP:32:VAL:HG12	28:DP:33:GLU:O	2.15	0.47
28:DP:98:TYR:C	28:DP:100:ARG:H	2.18	0.47
44:DQ:23:TYR:CD2	44:DQ:23:TYR:N	2.81	0.47
45:DS:69:LEU:HD23	45:DS:107:VAL:HG21	1.97	0.47
45:DS:73:LYS:HE3	45:DS:74:ILE:H	1.80	0.47
50:DT:40:LYS:C	50:DT:43:ILE:HG22	2.35	0.47
50:DT:40:LYS:HA	50:DT:43:ILE:HG22	1.95	0.47
35:DV:1:MET:O	35:DV:2:PHE:HB2	2.14	0.47
52:DW:50:VAL:HG23	52:DW:61:LYS:CE	2.44	0.47
1:AA:1009:U:H2'	1:AA:1010:U:C6	2.50	0.47
1:AA:1179:A:H2'	1:AA:1180:A:O4'	2.14	0.47
1:AA:131:A:H2'	1:AA:132:C:H6	1.78	0.47
1:AA:1358:U:H3'	1:AA:1359:C:C6	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1368:A:H5''	21:AN:100:TRP:CH2	2.50	0.47
1:AA:597:G:H2'	1:AA:598:U:H5'	1.95	0.47
1:AA:832:G:O2'	1:AA:833:G:H5'	2.15	0.47
18:AB:15:PHE:HD1	18:AB:16:GLY:N	2.13	0.47
18:AB:202:ASN:HB3	18:AB:208:ALA:HB3	1.96	0.47
2:AC:63:ILE:HB	2:AC:98:ALA:HB2	1.96	0.47
3:AD:176:LYS:HD3	3:AD:176:LYS:H	1.80	0.47
3:AD:178:GLU:HB3	3:AD:179:GLY:H	1.57	0.47
8:AI:20:ILE:HD13	8:AI:85:ALA:CB	2.39	0.47
9:AJ:52:LEU:HA	9:AJ:62:ARG:H	1.79	0.47
10:AK:22:ILE:HD13	10:AK:95:THR:HG21	1.96	0.47
12:AM:44:ILE:O	12:AM:47:LEU:HB2	2.15	0.47
21:AN:12:ARG:HD3	21:AN:58:ARG:O	2.15	0.47
16:AS:18:VAL:O	16:AS:22:VAL:HG23	2.15	0.47
17:AT:8:LYS:HE3	17:AT:12:GLN:OE1	2.14	0.47
23:BB:1124:G:N3	32:B4:38:GLY:O	2.48	0.47
22:BA:106:G:H2'	22:BA:107:G:O4'	2.14	0.47
22:BA:59:A:H2'	22:BA:60:C:H6	1.78	0.47
23:BB:1081:U:O2'	24:BI:118:GLY:HA2	2.15	0.47
23:BB:70:G:H2'	23:BB:113:U:O2'	2.14	0.47
23:BB:1404:C:H2'	23:BB:1405:U:C6	2.50	0.47
23:BB:1495:A:H2'	23:BB:1496:A:C8	2.49	0.47
23:BB:2026:U:H2'	23:BB:2027:G:C8	2.50	0.47
23:BB:2329:U:H2'	23:BB:2330:G:C8	2.50	0.47
23:BB:242:G:H8	34:B3:3:ILE:O	1.97	0.47
23:BB:413:C:H2'	23:BB:414:C:C6	2.50	0.47
23:BB:661:A:O2'	37:BL:13:LYS:HA	2.15	0.47
23:BB:664:G:H2'	23:BB:665:U:C6	2.49	0.47
23:BB:920:A:H2'	23:BB:921:C:C6	2.49	0.47
25:BC:123:ILE:HG13	25:BC:123:ILE:O	2.15	0.47
25:BC:149:LYS:HG2	25:BC:152:GLN:OE1	2.15	0.47
29:BE:101:TYR:O	29:BE:104:ALA:HB3	2.15	0.47
47:BF:113:PHE:HZ	47:BF:175:PRO:HB2	1.79	0.47
40:BH:5:LEU:HD13	40:BH:12:LEU:C	2.35	0.47
23:BB:1132:U:O2	41:BJ:75:TYR:HB2	2.15	0.47
41:BJ:93:ILE:CA	41:BJ:97:PRO:HG3	2.45	0.47
27:BK:108:ARG:HG3	27:BK:108:ARG:O	2.14	0.47
37:BL:110:VAL:HB	37:BL:127:VAL:CG2	2.44	0.47
37:BL:9:ALA:HB3	37:BL:12:SER:OG	2.14	0.47
31:B0:42:ILE:HD11	42:BN:98:LEU:HD12	1.97	0.47
43:BO:39:VAL:CG1	43:BO:48:LEU:HD12	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:87:LEU:HB2	50:BT:91:GLN:HG2	1.97	0.47
46:BU:78:LYS:HE3	46:BU:79:ALA:N	2.29	0.47
35:BV:42:LEU:HD23	35:BV:42:LEU:N	2.30	0.47
35:BV:60:VAL:HG12	35:BV:61:LEU:H	1.78	0.47
1:CA:1244:G:H2'	1:CA:1245:C:H6	1.76	0.47
1:CA:221:C:O2'	1:CA:222:C:H5'	2.14	0.47
1:CA:398:U:H2'	1:CA:399:G:H8	1.79	0.47
1:CA:783:C:O2'	1:CA:784:A:H5'	2.14	0.47
1:CA:857:C:H2'	1:CA:858:G:O4'	2.15	0.47
18:CB:30:ILE:HD11	18:CB:188:THR:HG22	1.96	0.47
18:CB:218:ALA:HA	18:CB:221:ARG:HH12	1.80	0.47
8:CI:79:ARG:O	8:CI:83:THR:HG22	2.15	0.47
1:CA:796:C:O3'	10:CK:126:ARG:NH2	2.47	0.47
1:CA:1227:A:H4'	12:CM:113:LYS:HE2	1.97	0.47
13:CP:68:SER:HB3	13:CP:71:VAL:HG12	1.96	0.47
34:D3:56:LEU:O	34:D3:59:ALA:HB3	2.15	0.47
23:DB:1292:G:O2'	23:DB:1293:C:H5'	2.15	0.47
23:DB:1327:A:H2'	23:DB:1328:A:H5'	1.96	0.47
23:DB:159:G:O2'	23:DB:160:A:H5''	2.15	0.47
23:DB:1714:U:H3'	23:DB:1715:G:H5'	1.97	0.47
23:DB:2078:C:H2'	23:DB:2079:U:H6	1.80	0.47
23:DB:2225:A:H4'	23:DB:2226:C:H6	1.80	0.47
23:DB:231:A:H3'	23:DB:232:G:C8	2.49	0.47
23:DB:2634:A:H2'	23:DB:2635:A:C8	2.50	0.47
23:DB:2649:C:H2'	23:DB:2650:U:C6	2.50	0.47
23:DB:2726:A:HO2'	23:DB:2727:A:C5'	2.28	0.47
23:DB:2783:U:H2'	23:DB:2784:U:C6	2.50	0.47
23:DB:404:A:H4'	23:DB:405:U:O5'	2.15	0.47
23:DB:437:U:H2'	23:DB:438:G:C8	2.50	0.47
23:DB:811:U:OP2	37:DL:20:GLY:HA2	2.15	0.47
23:DB:825:A:H2'	23:DB:826:U:O4'	2.14	0.47
23:DB:939:G:O2'	23:DB:940:G:H5'	2.14	0.47
29:DE:58:LYS:HD3	29:DE:58:LYS:N	2.29	0.47
29:DE:58:LYS:H	29:DE:58:LYS:HZ3	1.62	0.47
48:DG:102:ILE:O	48:DG:113:ASP:HA	2.15	0.47
48:DG:97:VAL:CG2	48:DG:124:CYS:HB2	2.44	0.47
48:DG:148:ARG:HA	48:DG:161:VAL:HB	1.96	0.47
24:DI:37:PHE:HB2	24:DI:66:PHE:CE2	2.50	0.47
37:DL:118:THR:O	37:DL:120:VAL:HG23	2.14	0.47
23:DB:833:A:H1'	37:DL:52:GLY:N	2.30	0.47
37:DL:81:ASP:HA	37:DL:84:LYS:CE	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:345:C:H5'	28:DP:38:ARG:NH2	2.29	0.47
49:DR:2:TYR:HB2	49:DR:42:ALA:N	2.30	0.47
45:DS:81:SER:HB3	45:DS:99:ARG:HB3	1.96	0.47
52:DW:28:GLU:O	52:DW:30:VAL:N	2.48	0.47
1:AA:1009:U:H1'	1:AA:1021:A:C2	2.50	0.47
1:AA:1218:C:H2'	1:AA:1219:A:H8	1.77	0.47
1:AA:1304:G:N1	1:AA:1305:G:N2	2.63	0.47
1:AA:1347:G:OP2	8:AI:108:ARG:HB2	2.15	0.47
1:AA:1382:C:H2'	1:AA:1383:C:H6	1.80	0.47
1:AA:67:C:H4'	1:AA:172:A:O4'	2.15	0.47
1:AA:425:G:O2'	1:AA:426:U:H5'	2.15	0.47
1:AA:552:U:H5'	11:AL:82:ARG:NH1	2.28	0.47
1:AA:634:C:H2'	1:AA:635:A:C8	2.50	0.47
1:AA:783:C:O2'	1:AA:784:A:H5'	2.14	0.47
1:AA:980:C:H2'	1:AA:981:U:H5'	1.97	0.47
18:AB:93:HIS:C	18:AB:94:ARG:HE	2.18	0.47
2:AC:184:ASN:ND2	2:AC:185:THR:H	2.13	0.47
3:AD:157:ALA:C	3:AD:159:GLU:H	2.18	0.47
5:AF:85:ILE:HG22	5:AF:86:ARG:N	2.30	0.47
5:AF:97:THR:HB	5:AF:98:GLU:OE1	2.15	0.47
9:AJ:7:ARG:C	9:AJ:8:ILE:HD12	2.35	0.47
13:AP:23:ASP:CG	13:AP:25:ARG:HE	2.18	0.47
36:B2:9:VAL:CG1	36:B2:10:LEU:H	2.26	0.47
22:BA:17:C:O2'	22:BA:18:G:H5'	2.15	0.47
23:BB:1102:C:H2'	23:BB:1103:A:H8	1.80	0.47
23:BB:1145:C:O2'	23:BB:1146:C:H5'	2.14	0.47
23:BB:1313:U:O2	23:BB:1313:U:C2'	2.62	0.47
23:BB:1486:U:H2'	23:BB:1487:U:H6	1.80	0.47
23:BB:1494:A:H2'	23:BB:1495:A:H8	1.80	0.47
23:BB:1521:G:H3'	23:BB:1522:A:H2'	1.97	0.47
23:BB:1869:G:N3	23:BB:1869:G:H2'	2.30	0.47
23:BB:1130:U:C2	23:BB:2025:C:H5''	2.50	0.47
23:BB:2260:C:O2'	23:BB:2261:C:H5'	2.15	0.47
23:BB:2300:C:H2'	23:BB:2301:C:H6	1.79	0.47
23:BB:231:A:H3'	23:BB:232:G:C8	2.49	0.47
23:BB:2590:A:H2'	23:BB:2591:C:H6	1.79	0.47
23:BB:2564:A:OP1	23:BB:2648:G:H4'	2.14	0.47
23:BB:281:C:H2'	23:BB:282:A:H8	1.80	0.47
23:BB:2876:G:H2'	23:BB:2877:G:O4'	2.15	0.47
23:BB:296:U:H2'	23:BB:297:G:H8	1.79	0.47
23:BB:476:G:H4'	23:BB:502:A:H61	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:727:A:O2'	23:BB:728:G:H5'	2.15	0.47
23:BB:814:C:H1'	23:BB:1225:G:N2	2.29	0.47
23:BB:825:A:H2'	23:BB:826:U:O4'	2.15	0.47
25:BC:110:LYS:O	25:BC:113:ASP:HB2	2.15	0.47
26:BD:121:THR:HG21	26:BD:127:PHE:CE1	2.50	0.47
23:BB:321:U:OP2	29:BE:130:LYS:HA	2.15	0.47
47:BF:107:VAL:HB	47:BF:108:PRO:HD3	1.97	0.47
47:BF:105:ILE:HG22	47:BF:109:ARG:HG2	1.96	0.47
48:BG:123:GLU:O	48:BG:125:PRO:HD3	2.15	0.47
40:BH:67:ALA:O	40:BH:70:GLU:HG2	2.15	0.47
27:BK:68:GLY:HA3	27:BK:78:ARG:HB3	1.96	0.47
37:BL:118:THR:O	37:BL:120:VAL:HG23	2.15	0.47
37:BL:69:ARG:HE	37:BL:69:ARG:HA	1.80	0.47
38:BM:10:ARG:HH21	38:BM:10:ARG:HG3	1.79	0.47
42:BN:28:LEU:CA	42:BN:34:ILE:HD11	2.45	0.47
42:BN:37:THR:HG22	42:BN:39:PRO:CD	2.35	0.47
23:BB:2821:A:OP2	42:BN:3:HIS:NE2	2.48	0.47
23:BB:1248:G:OP1	44:BQ:1:ALA:HB3	2.15	0.47
45:BS:33:LEU:HD23	45:BS:51:LEU:HD23	1.97	0.47
50:BT:45:ALA:O	50:BT:48:GLN:HB2	2.15	0.47
1:CA:1006:G:O2'	1:CA:1007:U:H5'	2.14	0.47
1:CA:1121:U:H2'	1:CA:1122:U:C6	2.50	0.47
1:CA:1296:C:C4'	1:CA:1302:C:N4	2.77	0.47
1:CA:1369:C:H2'	1:CA:1370:G:C8	2.50	0.47
1:CA:524:G:H2'	1:CA:525:C:H6	1.75	0.47
1:CA:93:U:O5'	1:CA:93:U:H6	1.97	0.47
2:CC:174:LEU:H	2:CC:174:LEU:HD12	1.80	0.47
3:CD:57:LYS:O	3:CD:61:ARG:HB2	2.15	0.47
6:CG:129:ASN:O	6:CG:130:LYS:HB2	2.15	0.47
10:CK:21:HIS:HA	10:CK:84:MET:O	2.15	0.47
10:CK:52:ARG:HB3	10:CK:52:ARG:NH1	2.30	0.47
12:CM:89:ARG:HH11	12:CM:94:LEU:HB3	1.79	0.47
21:CN:29:ILE:HG22	21:CN:30:ILE:H	1.80	0.47
16:CS:10:ILE:HG21	16:CS:40:PHE:CE1	2.50	0.47
19:CU:48:LYS:HG3	19:CU:49:ALA:N	2.28	0.47
36:D2:1:MET:HG2	36:D2:2:LYS:H	1.78	0.47
36:D2:31:LEU:CD2	36:D2:42:LEU:HD12	2.45	0.47
23:DB:1141:U:H5''	41:DJ:27:ARG:NH2	2.25	0.47
23:DB:1192:G:O2'	23:DB:1193:G:H5'	2.15	0.47
23:DB:1374:G:H2'	23:DB:1375:U:C6	2.49	0.47
23:DB:1404:C:H2'	23:DB:1405:U:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1469:A:H2'	23:DB:1470:A:C8	2.50	0.47
23:DB:1523:U:H5''	23:DB:1524:G:C8	2.50	0.47
23:DB:1682:G:H2'	23:DB:1683:U:C6	2.50	0.47
23:DB:1993:U:H4'	26:DD:133:THR:HG21	1.96	0.47
23:DB:2010:G:H2'	23:DB:2011:U:H6	1.80	0.47
23:DB:2018:G:O2'	23:DB:2019:A:H5'	2.15	0.47
23:DB:27:G:H1'	23:DB:513:A:N6	2.30	0.47
23:DB:2899:A:H2'	23:DB:2900:A:O4'	2.15	0.47
23:DB:393:C:O2'	23:DB:394:C:H5'	2.14	0.47
23:DB:494:G:O2'	23:DB:495:G:H5'	2.15	0.47
23:DB:527:C:O2	23:DB:527:C:O4'	2.31	0.47
23:DB:81:G:H2'	23:DB:82:U:O4'	2.15	0.47
25:DC:145:MET:HG3	25:DC:152:GLN:OE1	2.14	0.47
23:DB:2774:C:OP1	26:DD:169:ARG:HG3	2.14	0.47
29:DE:98:LYS:HE3	29:DE:99:LYS:HG2	1.97	0.47
48:DG:26:LYS:HB2	48:DG:32:LEU:HG	1.97	0.47
40:DH:128:HIS:O	40:DH:144:VAL:N	2.48	0.47
40:DH:39:ALA:HA	40:DH:44:ILE:HG21	1.97	0.47
40:DH:4:ILE:CG2	40:DH:16:GLY:HA2	2.44	0.47
38:DM:19:GLY:CA	38:DM:98:PRO:HD2	2.35	0.47
50:DT:45:ALA:O	50:DT:48:GLN:HB2	2.14	0.47
46:DU:13:LEU:CA	46:DU:18:LYS:HE3	2.44	0.47
23:DB:85:G:OP2	46:DU:6:ARG:HD3	2.15	0.47
23:DB:2264:C:H41	52:DW:11:ASN:HD22	1.63	0.47
1:AA:1144:G:H21	1:AA:1146:A:H62	1.61	0.47
1:AA:1314:C:H5	16:AS:5:LYS:HG3	1.80	0.47
1:AA:1392:G:H2'	1:AA:1393:U:O4'	2.14	0.47
1:AA:178:C:H2'	1:AA:179:A:H8	1.80	0.47
1:AA:386:C:C2'	1:AA:387:U:H5'	2.45	0.47
1:AA:450:G:N7	1:AA:481:G:O6	2.47	0.47
1:AA:470:C:H2'	1:AA:471:U:C6	2.50	0.47
1:AA:745:G:H5''	1:AA:851:G:O2'	2.15	0.47
18:AB:98:GLY:HA2	18:AB:101:THR:HG22	1.97	0.47
18:AB:221:ARG:HH11	18:AB:221:ARG:HB3	1.79	0.47
5:AF:64:VAL:HG12	5:AF:65:GLU:N	2.30	0.47
1:AA:598:U:H4'	7:AH:85:TYR:CG	2.50	0.47
9:AJ:5:ARG:NH1	9:AJ:5:ARG:HB2	2.30	0.47
34:B3:20:GLY:HA3	34:B3:48:MET:HE1	1.96	0.47
23:BB:1082:U:C2	23:BB:1086:A:N1	2.83	0.47
23:BB:1192:G:O2'	23:BB:1193:G:H5'	2.14	0.47
23:BB:1401:G:H2'	23:BB:1402:U:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:140:C:O5'	23:BB:141:G:N2	2.45	0.47
23:BB:1432:G:O2'	23:BB:1433:A:H5'	2.14	0.47
23:BB:1690:A:H2'	23:BB:1691:C:O4'	2.15	0.47
23:BB:2206:C:O2'	23:BB:2207:C:H5'	2.15	0.47
23:BB:2340:A:H2'	23:BB:2341:G:C8	2.49	0.47
23:BB:2467:C:H1'	38:BM:122:ALA:HB1	1.97	0.47
23:BB:2758:A:C2'	23:BB:2759:G:H5'	2.45	0.47
23:BB:463:G:N2	23:BB:466:A:OP2	2.44	0.47
23:BB:864:G:O2'	23:BB:865:C:H5'	2.15	0.47
26:BD:125:TRP:CE2	26:BD:160:LYS:HB3	2.50	0.47
29:BE:128:ALA:O	29:BE:130:LYS:N	2.45	0.47
48:BG:106:LEU:HD13	48:BG:151:ARG:HB2	1.97	0.47
48:BG:58:ALA:C	48:BG:60:GLY:H	2.18	0.47
41:BJ:55:ILE:CG2	41:BJ:123:LYS:HB2	2.41	0.47
22:BA:9:G:P	43:BO:25:ARG:HH12	2.38	0.47
28:BP:31:VAL:HG11	28:BP:38:ARG:HG2	1.95	0.47
44:BQ:96:ASP:C	44:BQ:98:ALA:H	2.17	0.47
30:BY:6:ILE:H	30:BY:6:ILE:CD1	2.25	0.47
1:CA:105:G:H2'	1:CA:106:C:H6	1.78	0.47
1:CA:1479:C:H2'	1:CA:1480:A:H8	1.79	0.47
1:CA:378:G:O2'	1:CA:379:C:H5'	2.15	0.47
1:CA:470:C:H2'	1:CA:471:U:C6	2.50	0.47
18:CB:44:LYS:O	18:CB:48:MET:HG2	2.15	0.47
1:CA:406:G:N2	3:CD:115:GLN:HE22	2.07	0.47
3:CD:28:ASP:CB	3:CD:33:ILE:HG21	2.45	0.47
33:D1:4:ILE:HB	33:D1:27:ARG:NH1	2.30	0.47
23:DB:1430:G:H2'	23:DB:1431:A:H8	1.79	0.47
23:DB:1495:A:H2'	23:DB:1496:A:C8	2.50	0.47
23:DB:1428:C:H2'	23:DB:1569:A:OP2	2.14	0.47
23:DB:1607:C:H5'	23:DB:1608:A:H5'	1.95	0.47
23:DB:1635:A:H2'	23:DB:1636:U:H5'	1.96	0.47
23:DB:1841:U:H2'	23:DB:1842:G:H8	1.78	0.47
23:DB:1131:G:N2	23:DB:2024:G:H21	2.13	0.47
23:DB:2147:A:H4'	23:DB:2148:G:H8	1.79	0.47
23:DB:225:C:H2'	23:DB:226:A:O4'	2.14	0.47
23:DB:2340:A:H2'	23:DB:2341:G:C8	2.50	0.47
23:DB:2722:G:H2'	23:DB:2723:C:O4'	2.14	0.47
23:DB:2805:C:O2'	23:DB:2806:C:H5'	2.15	0.47
23:DB:2867:G:N3	23:DB:2867:G:C2'	2.78	0.47
23:DB:944:C:H5'	23:DB:945:A:C5'	2.45	0.47
25:DC:199:HIS:C	25:DC:201:LEU:H	2.19	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:222:THR:HA	25:DC:231:HIS:O	2.15	0.47
47:DF:168:LEU:O	47:DF:169:LEU:HB2	2.15	0.47
48:DG:54:ARG:HE	48:DG:54:ARG:HA	1.80	0.47
24:DI:69:VAL:HG23	24:DI:69:VAL:O	2.14	0.47
38:DM:120:ALA:O	38:DM:123:LYS:HG2	2.15	0.47
28:DP:47:ILE:HD11	28:DP:59:THR:HG22	1.97	0.47
44:DQ:93:ILE:HG23	44:DQ:94:LEU:N	2.30	0.47
50:DT:58:VAL:O	50:DT:58:VAL:HG13	2.15	0.47
46:DU:41:VAL:HG22	46:DU:60:LYS:O	2.15	0.47
46:DU:86:PHE:CE1	46:DU:90:LYS:HB2	2.50	0.47
35:DV:75:GLN:HB3	35:DV:76:ASP:H	1.43	0.47
51:DZ:55:GLY:O	51:DZ:58:VAL:HB	2.15	0.47
1:AA:946:A:O2'	1:AA:1333:A:N3	2.46	0.46
1:AA:1390:U:H2'	1:AA:1391:U:O4'	2.15	0.46
1:AA:191:G:H2'	1:AA:192:A:C8	2.49	0.46
1:AA:45:G:O2'	1:AA:46:G:H5'	2.15	0.46
1:AA:792:A:H4'	1:AA:793:U:O5'	2.15	0.46
18:AB:46:VAL:HG13	18:AB:47:PRO:CD	2.39	0.46
2:AC:146:LYS:HE3	2:AC:202:PHE:HE2	1.80	0.46
6:AG:71:THR:HG22	6:AG:141:HIS:HE1	1.79	0.46
7:AH:54:THR:HG23	7:AH:55:LYS:N	2.30	0.46
8:AI:6:TYR:CG	8:AI:7:GLY:N	2.83	0.46
14:AQ:7:LEU:HD22	14:AQ:24:ILE:HG21	1.97	0.46
16:AS:10:ILE:HD12	16:AS:10:ILE:N	2.30	0.46
17:AT:70:LYS:HA	17:AT:73:ARG:CZ	2.45	0.46
31:B0:46:GLY:O	31:B0:53:VAL:N	2.48	0.46
22:BA:85:G:H2'	22:BA:86:G:C8	2.51	0.46
23:BB:1241:A:N3	23:BB:1241:A:H5'	2.30	0.46
23:BB:1446:C:H2'	23:BB:1447:C:C6	2.50	0.46
23:BB:1665:A:H2'	23:BB:1666:G:O4'	2.15	0.46
23:BB:1706:C:H2'	23:BB:1757:A:OP2	2.15	0.46
23:BB:1714:U:H3'	23:BB:1715:G:H5'	1.97	0.46
23:BB:2028:U:H2'	23:BB:2029:G:C8	2.50	0.46
23:BB:2600:A:O2'	23:BB:2601:C:H5'	2.15	0.46
23:BB:2803:G:H2'	23:BB:2804:U:C6	2.51	0.46
23:BB:2852:G:H2'	23:BB:2853:C:H6	1.78	0.46
23:BB:564:C:H1'	44:BQ:36:GLN:OE1	2.15	0.46
23:BB:630:G:N2	23:BB:632:A:H3'	2.30	0.46
23:BB:672:C:O2'	23:BB:673:C:H5'	2.14	0.46
23:BB:852:U:H2'	23:BB:853:C:C6	2.50	0.46
29:BE:189:THR:O	29:BE:193:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:67:ARG:NH1	29:BE:70:SER:OG	2.48	0.46
47:BF:169:LEU:HB3	47:BF:174:PHE:CD1	2.50	0.46
27:BK:60:ALA:CB	27:BK:86:LEU:HA	2.44	0.46
37:BL:78:ARG:NH2	37:BL:113:ALA:HB1	2.30	0.46
42:BN:90:ARG:HG2	42:BN:94:TYR:HD1	1.79	0.46
44:BQ:96:ASP:C	44:BQ:98:ALA:N	2.69	0.46
49:BR:22:LEU:HD23	49:BR:22:LEU:N	2.30	0.46
45:BS:4:ILE:HG21	45:BS:106:VAL:HG22	1.97	0.46
46:BU:10:VAL:HG22	46:BU:35:VAL:HG11	1.97	0.46
1:CA:1021:A:H2'	1:CA:1022:A:H8	1.80	0.46
1:CA:177:G:P	17:CT:59:ARG:HE	2.39	0.46
1:CA:426:U:H2'	1:CA:427:U:C6	2.50	0.46
1:CA:469:C:O2'	1:CA:470:C:H5'	2.15	0.46
1:CA:560:A:H4'	1:CA:561:U:C5'	2.36	0.46
1:CA:778:G:H2'	1:CA:779:C:H6	1.79	0.46
18:CB:211:LEU:O	18:CB:215:ALA:HB2	2.15	0.46
3:CD:90:LEU:N	3:CD:90:LEU:HD22	2.30	0.46
4:CE:113:VAL:O	4:CE:114:LEU:C	2.54	0.46
4:CE:36:THR:HB	4:CE:63:MET:HE2	1.96	0.46
5:CF:56:LYS:N	5:CF:56:LYS:HE3	2.29	0.46
7:CH:55:LYS:CE	7:CH:55:LYS:HA	2.45	0.46
8:CI:112:ARG:HH22	9:CJ:64:GLN:HE22	1.61	0.46
8:CI:94:ARG:NE	8:CI:97:LEU:HD12	2.29	0.46
34:D3:7:ARG:HG3	34:D3:7:ARG:HH11	1.79	0.46
23:DB:1056:G:H5''	23:DB:1057:A:H5'	1.97	0.46
23:DB:1279:G:H4'	42:DN:31:HIS:HD2	1.79	0.46
23:DB:1313:U:O2	23:DB:1313:U:C2'	2.63	0.46
23:DB:2097:A:H2'	23:DB:2098:U:C6	2.49	0.46
23:DB:381:G:H5''	51:DZ:16:ASN:ND2	2.30	0.46
23:DB:972:A:OP2	23:DB:974:G:H5''	2.15	0.46
25:DC:188:ARG:HG2	25:DC:188:ARG:HH21	1.79	0.46
25:DC:250:GLN:C	25:DC:252:LYS:H	2.18	0.46
29:DE:148:ILE:HG13	29:DE:167:VAL:CG2	2.41	0.46
29:DE:172:ALA:HB3	29:DE:195:GLN:NE2	2.30	0.46
29:DE:21:ARG:HG2	29:DE:107:SER:HB3	1.98	0.46
47:DF:116:LEU:HD23	47:DF:174:PHE:CE2	2.50	0.46
47:DF:7:TYR:OH	47:DF:29:ARG:HG3	2.16	0.46
48:DG:15:ASP:O	48:DG:16:VAL:HB	2.15	0.46
48:DG:58:ALA:C	48:DG:60:GLY:H	2.18	0.46
40:DH:18:GLN:HE21	40:DH:39:ALA:HB1	1.80	0.46
40:DH:74:ALA:O	40:DH:76:GLU:HG3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:DI:10:LEU:HD12	24:DI:10:LEU:O	2.14	0.46
24:DI:27:LEU:HD23	24:DI:27:LEU:N	2.22	0.46
23:DB:632:A:H1'	37:DL:66:PHE:HE2	1.81	0.46
37:DL:95:LEU:O	37:DL:100:ILE:HG22	2.16	0.46
38:DM:100:LYS:CD	38:DM:101:VAL:H	2.27	0.46
38:DM:64:TRP:C	38:DM:65:ILE:HG13	2.35	0.46
43:DO:75:GLY:O	43:DO:78:VAL:HG23	2.15	0.46
28:DP:100:ARG:HB3	28:DP:101:GLU:H	1.44	0.46
45:DS:76:VAL:HG12	45:DS:103:ILE:HA	1.97	0.46
23:DB:2352:A:C2	52:DW:29:SER:HB3	2.50	0.46
30:DY:6:ILE:HG21	30:DY:47:ILE:HD11	1.96	0.46
1:AA:1142:G:C2	1:AA:1143:G:H1'	2.50	0.46
1:AA:177:G:P	17:AT:59:ARG:HE	2.38	0.46
1:AA:956:U:O2'	1:AA:957:U:H5'	2.15	0.46
4:AE:40:ASP:OD2	4:AE:42:ASN:HB3	2.14	0.46
10:AK:86:LYS:HB2	10:AK:113:THR:HA	1.98	0.46
12:AM:12:LYS:O	12:AM:43:LYS:HG2	2.14	0.46
12:AM:67:ASP:O	12:AM:71:GLU:HG3	2.15	0.46
12:AM:76:ILE:HG22	12:AM:80:MET:HE3	1.96	0.46
23:BB:1100:C:H2'	23:BB:1101:U:C6	2.50	0.46
23:BB:1295:C:H2'	23:BB:1296:G:C8	2.50	0.46
23:BB:17:G:H2'	23:BB:18:U:H6	1.80	0.46
23:BB:2312:U:C4'	47:BF:84:ILE:HG21	2.44	0.46
23:BB:2415:G:H2'	23:BB:2416:C:H6	1.80	0.46
23:BB:340:A:H2'	23:BB:341:C:H5'	1.96	0.46
23:BB:521:U:H2'	23:BB:522:A:C8	2.50	0.46
23:BB:574:A:H4'	23:BB:575:A:C5'	2.45	0.46
23:BB:660:C:H2'	23:BB:661:A:H8	1.80	0.46
23:BB:942:G:H2'	23:BB:943:A:H8	1.79	0.46
23:BB:962:G:H21	23:BB:2250:G:H22	1.62	0.46
25:BC:68:ARG:HB3	25:BC:128:THR:OG1	2.15	0.46
23:BB:2060:A:C2'	29:BE:63:LYS:HE2	2.39	0.46
47:BF:21:TYR:CD2	47:BF:27:VAL:HG12	2.51	0.46
47:BF:24:VAL:C	47:BF:26:GLN:H	2.19	0.46
48:BG:134:GLY:HA3	48:BG:140:ILE:HG21	1.96	0.46
41:BJ:45:THR:HG23	41:BJ:45:THR:O	2.15	0.46
23:BB:2548:U:H1'	27:BK:23:LYS:HZ1	1.80	0.46
42:BN:12:ARG:HG3	42:BN:13:ASN:N	2.31	0.46
43:BO:14:ALA:C	43:BO:16:ARG:H	2.18	0.46
28:BP:47:ILE:HA	28:BP:96:LEU:HB2	1.96	0.46
46:BU:34:ILE:HG12	46:BU:63:ALA:CB	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:176:C:H3'	1:CA:177:G:H21	1.80	0.46
1:CA:635:A:H2'	1:CA:636:U:C6	2.50	0.46
1:CA:846:G:H2'	1:CA:846:G:N3	2.30	0.46
4:CE:21:SER:HB3	4:CE:28:ARG:NH2	2.30	0.46
4:CE:61:LYS:HG3	4:CE:62:ALA:N	2.29	0.46
6:CG:27:ASN:O	6:CG:30:MET:HB3	2.15	0.46
1:CA:598:U:H4'	7:CH:85:TYR:CG	2.50	0.46
10:CK:69:CYS:SG	10:CK:70:ALA:N	2.88	0.46
11:CL:15:VAL:O	11:CL:16:ALA:C	2.53	0.46
11:CL:32:VAL:HB	11:CL:55:ARG:HB3	1.97	0.46
12:CM:15:VAL:HG13	12:CM:33:LEU:HD11	1.97	0.46
16:CS:52:ASN:C	16:CS:76:THR:HG23	2.35	0.46
34:D3:21:PHE:CE1	34:D3:58:ILE:HG23	2.50	0.46
22:DA:109:A:H2'	22:DA:110:C:C6	2.50	0.46
23:DB:1242:U:H2'	23:DB:1243:C:C6	2.49	0.46
23:DB:2026:U:H2'	23:DB:2027:G:C8	2.51	0.46
23:DB:2041:U:H2'	23:DB:2042:A:H8	1.79	0.46
23:DB:2109:U:H6	23:DB:2109:U:OP1	1.99	0.46
23:DB:2479:U:H2'	23:DB:2480:C:H5'	1.97	0.46
23:DB:2648:G:H2'	23:DB:2649:C:C6	2.50	0.46
23:DB:2803:G:H2'	23:DB:2804:U:C6	2.50	0.46
23:DB:2849:U:H4'	23:DB:2850:A:C5'	2.44	0.46
23:DB:2901:C:H2'	23:DB:2902:C:O4'	2.15	0.46
23:DB:402:A:O2'	23:DB:403:U:H5'	2.15	0.46
23:DB:562:U:C4	23:DB:2036:C:O4'	2.68	0.46
23:DB:686:U:H1'	36:D2:6:GLN:O	2.16	0.46
23:DB:747:U:P	45:DS:90:LYS:HZ2	2.38	0.46
25:DC:163:ILE:HG22	25:DC:164:VAL:N	2.30	0.46
23:DB:1654:A:O2'	26:DD:118:PHE:HB3	2.15	0.46
29:DE:114:ARG:HG3	29:DE:114:ARG:HH11	1.80	0.46
29:DE:138:LEU:HD22	29:DE:143:LEU:HB2	1.97	0.46
47:DF:105:ILE:HG22	47:DF:109:ARG:HG2	1.96	0.46
47:DF:4:HIS:O	47:DF:7:TYR:HB3	2.15	0.46
40:DH:77:THR:HG22	40:DH:143:ILE:CB	2.42	0.46
41:DJ:56:VAL:C	41:DJ:57:LEU:HD12	2.36	0.46
41:DJ:19:ASP:CB	41:DJ:57:LEU:HB2	2.45	0.46
27:DK:70:ARG:CB	27:DK:76:VAL:HG22	2.43	0.46
42:DN:32:GLU:HG2	42:DN:115:LEU:HD12	1.97	0.46
42:DN:29:VAL:HG13	42:DN:83:LEU:HD22	1.97	0.46
28:DP:103:THR:HG22	28:DP:104:GLY:N	2.26	0.46
28:DP:19:PHE:HE2	28:DP:83:ILE:HD11	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DR:19:THR:HA	49:DR:97:LYS:HA	1.97	0.46
45:DS:52:GLU:HA	45:DS:55:ILE:HG22	1.97	0.46
46:DU:35:VAL:HB	46:DU:38:ILE:HG21	1.97	0.46
46:DU:73:ASN:OD1	46:DU:75:ALA:HB3	2.15	0.46
1:AA:1081:A:O2'	1:AA:1082:A:H5'	2.16	0.46
1:AA:1360:A:N6	1:AA:1361:G:N2	2.63	0.46
1:AA:1395:C:O2'	1:AA:1396:A:H5'	2.15	0.46
1:AA:24:U:O2'	1:AA:25:C:H5'	2.15	0.46
1:AA:45:G:H2'	1:AA:46:G:C8	2.48	0.46
1:AA:475:C:O2'	1:AA:476:U:H5'	2.16	0.46
1:AA:764:C:N4	1:AA:812:G:H1	2.13	0.46
18:AB:160:LEU:O	18:AB:182:VAL:HG13	2.15	0.46
3:AD:111:ALA:O	3:AD:114:ARG:HB2	2.16	0.46
4:AE:89:THR:HB	4:AE:134:ASN:ND2	2.29	0.46
4:AE:24:VAL:HG23	4:AE:26:GLY:H	1.79	0.46
6:AG:22:LEU:O	6:AG:25:PHE:HB3	2.16	0.46
8:AI:21:LYS:O	8:AI:60:LEU:HB2	2.15	0.46
1:AA:1308:U:H5''	12:AM:96:VAL:CG2	2.45	0.46
21:AN:1:ALA:O	21:AN:6:LYS:HG3	2.16	0.46
14:AQ:26:ARG:N	14:AQ:39:ARG:O	2.48	0.46
10:AK:111:ASP:HB3	19:AU:19:LYS:HE2	1.97	0.46
23:BB:1386:C:H2'	23:BB:1387:A:C8	2.51	0.46
23:BB:2559:C:O2'	23:BB:2560:A:H5'	2.15	0.46
23:BB:286:U:O2'	23:BB:287:G:H5'	2.15	0.46
23:BB:475:C:H4'	23:BB:509:C:H2'	1.97	0.46
23:BB:81:G:H2'	23:BB:82:U:O4'	2.16	0.46
25:BC:232:GLY:N	25:BC:241:LYS:NZ	2.64	0.46
27:BK:79:PHE:HD2	28:BP:69:VAL:HG12	1.80	0.46
42:BN:62:ASN:N	42:BN:62:ASN:ND2	2.62	0.46
42:BN:9:GLN:HA	42:BN:17:ARG:CD	2.45	0.46
43:BO:105:ALA:O	43:BO:106:LEU:HG	2.15	0.46
28:BP:19:PHE:O	28:BP:20:ARG:HB2	2.14	0.46
28:BP:47:ILE:HG13	28:BP:48:ALA:H	1.80	0.46
44:BQ:59:LEU:O	44:BQ:62:ALA:HB3	2.14	0.46
44:BQ:86:SER:HB3	49:BR:50:GLY:O	2.16	0.46
49:BR:34:GLU:OE1	49:BR:60:LYS:HG2	2.15	0.46
46:BU:26:ASN:HD22	46:BU:26:ASN:H	1.61	0.46
46:BU:3:LYS:HD3	46:BU:82:VAL:CG2	2.46	0.46
52:BW:13:ARG:HA	52:BW:13:ARG:HH11	1.79	0.46
1:CA:112:G:O2'	1:CA:113:G:H5'	2.15	0.46
1:CA:1473:G:H2'	1:CA:1474:U:O4'	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1499:A:O2'	1:CA:1500:A:H5'	2.15	0.46
1:CA:59:A:H1'	1:CA:354:G:N2	2.31	0.46
1:CA:525:C:H2'	1:CA:526:C:C6	2.49	0.46
1:CA:724:G:H2'	1:CA:725:G:C8	2.45	0.46
1:CA:77:A:H2'	1:CA:78:A:H8	1.80	0.46
1:CA:745:G:H5''	1:CA:851:G:O2'	2.14	0.46
1:CA:94:G:H4'	1:CA:95:C:H5''	1.98	0.46
18:CB:17:HIS:HB2	18:CB:202:ASN:OD1	2.15	0.46
18:CB:53:LEU:HD22	18:CB:212:TYR:CZ	2.50	0.46
3:CD:153:ARG:HG3	3:CD:154:VAL:N	2.26	0.46
3:CD:48:SER:O	3:CD:49:ASP:C	2.54	0.46
4:CE:17:VAL:HG23	4:CE:33:THR:O	2.16	0.46
4:CE:24:VAL:HG23	4:CE:26:GLY:H	1.80	0.46
5:CF:92:THR:HG22	5:CF:94:HIS:N	2.17	0.46
6:CG:14:ASP:OD1	6:CG:17:PHE:N	2.48	0.46
1:CA:1129:C:H5''	8:CI:17:ARG:NH1	2.29	0.46
20:CO:7:ALA:HA	20:CO:10:LYS:HB3	1.97	0.46
1:CA:1320:C:C5	16:CS:36:ARG:HA	2.51	0.46
17:CT:47:GLN:HE21	17:CT:82:ILE:HD11	1.81	0.46
31:D0:39:ARG:O	31:D0:40:HIS:HB2	2.15	0.46
33:D1:8:ILE:HG23	33:D1:9:LYS:N	2.31	0.46
34:D3:31:ILE:HD11	34:D3:34:LYS:CD	2.41	0.46
34:D3:41:ARG:C	34:D3:43:LEU:H	2.17	0.46
23:DB:1542:U:H2'	23:DB:1543:G:O4'	2.15	0.46
23:DB:1700:A:H2'	23:DB:1701:A:H5'	1.97	0.46
23:DB:212:G:H2'	23:DB:213:A:C8	2.51	0.46
23:DB:2730:C:H2'	23:DB:2731:G:C8	2.49	0.46
23:DB:275:C:H2'	23:DB:276:U:C4'	2.32	0.46
23:DB:2835:A:H62	23:DB:2878:U:H2'	1.80	0.46
23:DB:416:U:H2'	23:DB:417:C:C6	2.50	0.46
23:DB:445:C:O2'	23:DB:446:G:H5'	2.15	0.46
23:DB:581:C:H5'	44:DQ:30:VAL:HG23	1.97	0.46
23:DB:586:A:H5'	29:DE:84:THR:OG1	2.15	0.46
23:DB:660:C:H2'	23:DB:661:A:C8	2.50	0.46
23:DB:969:G:H2'	23:DB:970:U:C6	2.50	0.46
25:DC:141:HIS:NE2	25:DC:194:VAL:HA	2.30	0.46
25:DC:169:ALA:O	25:DC:185:ALA:HB3	2.15	0.46
25:DC:166:ARG:HA	25:DC:171:VAL:HG22	1.98	0.46
25:DC:232:GLY:N	25:DC:241:LYS:NZ	2.63	0.46
25:DC:78:GLU:HB2	25:DC:92:LEU:HG	1.97	0.46
26:DD:14:ILE:HA	28:DP:11:GLN:OE1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DE:3:LEU:CB	29:DE:12:LEU:HB3	2.44	0.46
47:DF:26:GLN:O	47:DF:27:VAL:C	2.54	0.46
47:DF:41:GLU:CB	47:DF:48:LEU:HD11	2.46	0.46
48:DG:17:LYS:HE2	48:DG:19:ASN:CG	2.36	0.46
41:DJ:42:ALA:O	41:DJ:44:TYR:N	2.49	0.46
23:DB:2295:C:OP2	43:DO:10:ARG:HD3	2.15	0.46
23:DB:2845:U:O3'	28:DP:52:ARG:HD3	2.15	0.46
44:DQ:24:TYR:CD1	44:DQ:25:GLY:N	2.84	0.46
49:DR:34:GLU:HB3	49:DR:58:VAL:HG21	1.97	0.46
46:DU:24:VAL:HA	46:DU:35:VAL:HA	1.97	0.46
51:DZ:7:VAL:HG23	51:DZ:67:VAL:HG11	1.98	0.46
1:AA:1378:C:C4	1:AA:1379:G:H1'	2.50	0.46
1:AA:430:A:O2'	1:AA:431:A:H5'	2.16	0.46
1:AA:659:U:H2'	1:AA:660:C:C6	2.50	0.46
1:AA:852:G:H2'	1:AA:853:C:H6	1.79	0.46
1:AA:86:G:C2	1:AA:87:C:N4	2.83	0.46
18:AB:162:VAL:HG22	18:AB:163:ILE:N	2.29	0.46
18:AB:23:ASN:HD21	18:AB:25:LYS:CD	2.28	0.46
2:AC:195:ILE:HG22	2:AC:195:ILE:O	2.16	0.46
3:AD:108:ALA:H	3:AD:112:GLU:CD	2.19	0.46
6:AG:21:LEU:N	6:AG:21:LEU:HD23	2.30	0.46
8:AI:41:GLU:N	8:AI:44:ARG:NH1	2.63	0.46
10:AK:83:VAL:CG2	10:AK:109:ILE:HG12	2.45	0.46
11:AL:74:GLN:HB3	11:AL:75:GLU:H	1.48	0.46
20:AO:25:THR:HG21	20:AO:70:LEU:CD2	2.45	0.46
19:AU:48:LYS:HG3	19:AU:49:ALA:N	2.29	0.46
34:B3:28:LEU:O	34:B3:29:ARG:HB3	2.16	0.46
23:BB:1174:U:O4'	23:BB:1174:U:O2	2.34	0.46
23:BB:1178:C:H2'	23:BB:1179:G:H8	1.80	0.46
23:BB:118:A:OP2	23:BB:119:A:H2'	2.15	0.46
23:BB:136:G:OP2	23:BB:137:U:H5	1.99	0.46
23:BB:1559:U:H3'	23:BB:1560:G:H5'	1.97	0.46
23:BB:2080:A:H4'	51:BZ:19:SER:OG	2.16	0.46
23:BB:2282:G:H4'	23:BB:2283:C:O5'	2.15	0.46
23:BB:2526:G:H1'	32:B4:1:MET:N	2.31	0.46
23:BB:2538:C:H2'	23:BB:2539:C:H6	1.80	0.46
23:BB:2734:A:C2'	23:BB:2735:G:H5'	2.44	0.46
23:BB:296:U:O2'	23:BB:297:G:H5'	2.15	0.46
23:BB:55:G:H2'	23:BB:56:A:H8	1.81	0.46
23:BB:680:C:H2'	23:BB:681:G:H8	1.79	0.46
23:BB:1799:G:N2	25:BC:153:LEU:HD22	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:247:TRP:O	25:BC:249:VAL:HG22	2.15	0.46
25:BC:68:ARG:HH21	25:BC:190:THR:CG2	2.28	0.46
26:BD:47:ALA:HB2	26:BD:83:ARG:HD2	1.96	0.46
47:BF:21:TYR:HD2	47:BF:27:VAL:HG12	1.80	0.46
24:BI:19:PRO:HB2	24:BI:22:PRO:HD2	1.97	0.46
27:BK:99:ILE:HD13	27:BK:118:LEU:CD2	2.45	0.46
27:BK:71:ARG:NE	27:BK:72:PRO:HD3	2.30	0.46
38:BM:59:ARG:HE	38:BM:60:GLN:CG	2.10	0.46
49:BR:14:VAL:HG23	49:BR:18:GLN:HG3	1.96	0.46
50:BT:28:ASN:ND2	50:BT:29:THR:HG23	2.30	0.46
52:BW:23:LYS:CG	52:BW:24:ARG:N	2.77	0.46
52:BW:30:VAL:O	52:BW:30:VAL:HG13	2.16	0.46
1:CA:1508:A:H2'	1:CA:1509:C:H6	1.78	0.46
1:CA:415:A:OP1	1:CA:415:A:H4'	2.16	0.46
18:CB:98:GLY:O	18:CB:100:LEU:N	2.42	0.46
1:CA:1107:C:OP1	2:CC:171:ARG:HB2	2.16	0.46
2:CC:179:ALA:HB1	2:CC:202:PHE:HE1	1.80	0.46
3:CD:18:LEU:HD23	3:CD:20:LEU:HD11	1.97	0.46
4:CE:152:VAL:O	4:CE:156:ARG:HG2	2.16	0.46
6:CG:112:ASP:HB3	6:CG:113:LYS:H	1.48	0.46
9:CJ:46:LYS:O	9:CJ:48:ARG:HG3	2.16	0.46
9:CJ:92:LEU:HD13	9:CJ:92:LEU:N	2.30	0.46
10:CK:121:ARG:HH12	19:CU:37:TYR:HE2	1.63	0.46
12:CM:13:HIS:HB3	12:CM:41:ASP:HA	1.97	0.46
16:CS:68:HIS:HB3	16:CS:72:GLU:OE2	2.16	0.46
36:D2:13:ASN:N	36:D2:13:ASN:HD22	2.13	0.46
23:DB:1351:C:O2'	23:DB:1571:A:H1'	2.15	0.46
23:DB:1635:A:O2'	23:DB:1636:U:H5'	2.15	0.46
23:DB:1727:C:H2'	23:DB:1728:C:H6	1.80	0.46
23:DB:2137:U:H2'	23:DB:2138:G:C8	2.50	0.46
23:DB:2252:G:O2'	23:DB:2253:G:H5'	2.15	0.46
23:DB:2897:U:H2'	23:DB:2898:U:H6	1.81	0.46
23:DB:691:C:O2'	23:DB:692:C:H5'	2.15	0.46
23:DB:776:G:H4'	23:DB:777:G:O5'	2.16	0.46
23:DB:789:A:H4'	23:DB:790:U:OP2	2.15	0.46
25:DC:103:ILE:HG22	25:DC:104:LEU:N	2.31	0.46
25:DC:42:ARG:NH2	25:DC:48:ILE:HD11	2.30	0.46
26:DD:109:VAL:HG11	26:DD:193:VAL:CG1	2.45	0.46
47:DF:31:GLU:HG3	47:DF:32:LYS:HG2	1.96	0.46
48:DG:146:ASP:O	48:DG:150:TYR:HD1	1.98	0.46
48:DG:84:LYS:HG2	48:DG:85:LYS:N	2.25	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DO:100:HIS:C	43:DO:104:GLN:HB2	2.36	0.46
43:DO:17:LYS:O	43:DO:21:LEU:HB2	2.14	0.46
43:DO:37:ALA:HB1	43:DO:78:VAL:HG21	1.98	0.46
28:DP:47:ILE:HG22	28:DP:96:LEU:HB3	1.97	0.46
44:DQ:57:ARG:C	44:DQ:59:LEU:H	2.18	0.46
45:DS:57:ASN:HD22	45:DS:57:ASN:N	2.11	0.46
35:DV:42:LEU:HD12	35:DV:47:VAL:HG21	1.97	0.46
1:AA:1179:A:H2'	1:AA:1180:A:H8	1.80	0.46
1:AA:20:U:H2'	1:AA:21:G:O4'	2.14	0.46
1:AA:239:U:H5''	1:AA:239:U:H6	1.79	0.46
18:AB:86:CYS:SG	18:AB:87:ASP:N	2.88	0.46
2:AC:63:ILE:HG22	2:AC:65:VAL:CG2	2.46	0.46
5:AF:3:HIS:CD2	5:AF:92:THR:HG23	2.51	0.46
6:AG:111:GLY:HA2	6:AG:118:ARG:NH1	2.30	0.46
8:AI:19:PHE:O	8:AI:62:LEU:HA	2.15	0.46
12:AM:13:HIS:HB3	12:AM:40:GLU:O	2.15	0.46
9:AJ:51:VAL:HG23	21:AN:80:ARG:HB2	1.98	0.46
20:AO:83:GLU:C	20:AO:85:LEU:H	2.19	0.46
14:AQ:10:ARG:HA	14:AQ:56:ASP:O	2.15	0.46
15:AR:27:THR:O	15:AR:28:LEU:HD23	2.16	0.46
33:B1:8:ILE:HD12	33:B1:9:LYS:H	1.81	0.46
34:B3:40:LYS:O	34:B3:43:LEU:HB2	2.15	0.46
32:B4:27:CYS:HB3	32:B4:33:HIS:HB2	1.98	0.46
23:BB:1464:G:H2'	23:BB:1465:G:H8	1.80	0.46
23:BB:1717:A:H2'	23:BB:1718:G:O4'	2.15	0.46
23:BB:1867:G:O2'	23:BB:1868:C:H5'	2.16	0.46
23:BB:1914:C:H2'	23:BB:1915:U:H5''	1.96	0.46
23:BB:2105:U:H2'	23:BB:2106:U:C6	2.51	0.46
23:BB:2623:G:O2'	23:BB:2624:G:H5'	2.16	0.46
23:BB:264:C:HO2'	23:BB:429:A:H2	1.62	0.46
23:BB:2686:G:H2'	23:BB:2687:U:C6	2.51	0.46
23:BB:274:C:O5'	23:BB:274:C:H6	1.98	0.46
23:BB:418:C:H2'	23:BB:419:U:C6	2.51	0.46
23:BB:564:C:O2'	23:BB:565:C:H5'	2.15	0.46
23:BB:625:G:O2'	23:BB:626:A:H5'	2.16	0.46
25:BC:222:THR:HA	25:BC:231:HIS:O	2.15	0.46
25:BC:32:LEU:HD22	25:BC:63:ILE:HG21	1.97	0.46
29:BE:60:TRP:CE3	29:BE:62:GLN:HA	2.51	0.46
29:BE:58:LYS:O	29:BE:60:TRP:N	2.49	0.46
47:BF:147:ARG:HB3	47:BF:147:ARG:CZ	2.46	0.46
38:BM:130:PHE:O	38:BM:131:VAL:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BS:29:VAL:HG23	45:BS:70:LYS:HA	1.98	0.46
50:BT:68:LYS:HD3	50:BT:68:LYS:N	2.31	0.46
51:BZ:5:CYS:HB3	51:BZ:9:GLY:N	2.30	0.46
1:CA:1459:G:O2'	1:CA:1460:C:H5'	2.15	0.46
1:CA:263:A:OP1	17:CT:73:ARG:NH1	2.49	0.46
1:CA:295:C:H2'	1:CA:296:U:H6	1.81	0.46
1:CA:57:G:N2	1:CA:388:G:C6	2.84	0.46
1:CA:737:C:H2'	1:CA:738:C:H6	1.79	0.46
5:CF:51:ILE:O	5:CF:51:ILE:HG23	2.15	0.46
9:CJ:5:ARG:N	9:CJ:77:VAL:HA	2.30	0.46
12:CM:90:HIS:HA	12:CM:108:ARG:HH22	1.78	0.46
12:CM:78:ARG:HH12	16:CS:64:GLU:HG2	1.80	0.46
22:DA:112:G:O2'	22:DA:113:C:H5'	2.15	0.46
22:DA:37:C:H3'	22:DA:38:C:H6	1.81	0.46
23:DB:1040:A:O2'	23:DB:1041:G:H5'	2.15	0.46
23:DB:1093:G:O2'	23:DB:1094:U:H5'	2.15	0.46
23:DB:1745:A:O2'	23:DB:1746:A:H5'	2.16	0.46
23:DB:1846:G:N2	23:DB:1848:A:N6	2.64	0.46
23:DB:1859:U:H2'	23:DB:1860:G:H8	1.78	0.46
23:DB:577:G:OP1	23:DB:2502:G:H2'	2.15	0.46
23:DB:2756:U:H6	23:DB:2756:U:O5'	1.99	0.46
23:DB:406:G:O2'	23:DB:407:G:H5'	2.16	0.46
23:DB:457:A:N1	23:DB:470:A:H5''	2.30	0.46
23:DB:483:A:H5''	46:DU:46:LYS:HG3	1.96	0.46
47:DF:3:LEU:HD12	47:DF:96:TRP:CD1	2.50	0.46
24:DI:52:LEU:HD13	24:DI:81:LYS:NZ	2.31	0.46
27:DK:99:ILE:HD13	27:DK:118:LEU:CD2	2.45	0.46
27:DK:72:PRO:O	27:DK:74:GLY:N	2.48	0.46
43:DO:31:THR:HG23	43:DO:34:HIS:O	2.15	0.46
50:DT:74:ILE:HG13	50:DT:75:GLY:N	2.31	0.46
35:DV:42:LEU:HB2	35:DV:47:VAL:HG21	1.98	0.46
39:DX:22:LEU:HG	39:DX:23:ARG:HG2	1.98	0.46
39:DX:7:ARG:HA	39:DX:7:ARG:HE	1.79	0.46
1:AA:105:G:H2'	1:AA:106:C:H6	1.78	0.46
1:AA:1249:C:H4'	8:AI:37:TYR:OH	2.16	0.46
1:AA:1505:G:H5''	1:AA:1506:U:O5'	2.15	0.46
1:AA:518:C:H2'	1:AA:530:G:N7	2.31	0.46
1:AA:692:U:C2	1:AA:694:A:H5''	2.50	0.46
1:AA:897:C:O2'	1:AA:898:G:H5'	2.16	0.46
1:AA:8:A:H1'	4:AE:107:GLY:HA2	1.97	0.46
1:AA:93:U:H2'	1:AA:94:G:H4'	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:160:LEU:O	18:AB:182:VAL:HA	2.16	0.46
18:AB:75:ALA:O	18:AB:76:SER:C	2.53	0.46
2:AC:119:ILE:HG23	2:AC:120:THR:N	2.30	0.46
3:AD:146:GLU:HA	3:AD:149:LYS:CG	2.45	0.46
4:AE:132:PRO:HG2	4:AE:133:ILE:H	1.80	0.46
7:AH:44:PHE:CE2	7:AH:100:ILE:HG12	2.50	0.46
8:AI:27:ILE:HG21	8:AI:34:LEU:HD13	1.95	0.46
8:AI:39:GLY:O	8:AI:40:ARG:HB2	2.15	0.46
11:AL:15:VAL:O	11:AL:16:ALA:C	2.53	0.46
20:AO:69:TYR:HA	20:AO:72:ARG:NH2	2.31	0.46
16:AS:39:ILE:HG22	16:AS:40:PHE:N	2.31	0.46
16:AS:38:THR:HG23	16:AS:69:LYS:CE	2.46	0.46
34:B3:60:CYS:C	34:B3:62:PRO:HD3	2.36	0.46
22:BA:109:A:H2'	22:BA:110:C:C6	2.51	0.46
23:BB:1635:A:O2'	23:BB:1636:U:H5'	2.14	0.46
23:BB:1654:A:H2'	23:BB:1655:A:C8	2.51	0.46
23:BB:1767:G:O2'	23:BB:1768:C:H5'	2.16	0.46
23:BB:1877:A:H2'	23:BB:1878:G:C8	2.50	0.46
23:BB:2471:A:O2'	23:BB:2472:G:O5'	2.34	0.46
23:BB:2488:G:H2'	23:BB:2489:U:C6	2.50	0.46
23:BB:341:C:H2'	23:BB:342:A:C8	2.51	0.46
23:BB:969:G:H2'	23:BB:970:U:C6	2.50	0.46
26:BD:56:LYS:HG3	26:BD:58:ASN:HB3	1.97	0.46
29:BE:21:ARG:HG2	29:BE:107:SER:HB3	1.97	0.46
29:BE:108:ILE:O	29:BE:108:ILE:HD13	2.16	0.46
23:BB:321:U:OP2	29:BE:130:LYS:HD3	2.15	0.46
29:BE:188:MET:HG2	29:BE:193:VAL:CG2	2.46	0.46
40:BH:4:ILE:CG2	40:BH:16:GLY:HA2	2.45	0.46
41:BJ:64:VAL:CG2	41:BJ:68:LYS:HB2	2.45	0.46
27:BK:118:LEU:O	27:BK:120:PRO:HD2	2.15	0.46
23:BB:2547:A:H4'	27:BK:29:HIS:CE1	2.50	0.46
42:BN:72:ASP:O	42:BN:76:VAL:HG13	2.14	0.46
42:BN:96:ARG:HD3	42:BN:96:ARG:N	2.31	0.46
28:BP:19:PHE:CE1	28:BP:46:VAL:HG11	2.50	0.46
49:BR:14:VAL:HG21	49:BR:98:ILE:CD1	2.46	0.46
45:BS:24:ILE:HG12	45:BS:36:LEU:HD21	1.98	0.46
50:BT:58:VAL:O	50:BT:58:VAL:HG13	2.16	0.46
50:BT:55:VAL:CA	50:BT:87:LEU:HA	2.30	0.46
51:BZ:19:SER:OG	51:BZ:23:ASN:HB2	2.16	0.46
51:BZ:53:ALA:O	51:BZ:54:LYS:HB3	2.16	0.46
1:CA:1326:U:H2'	1:CA:1327:C:H6	1.77	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1514:G:O2'	1:CA:1515:G:H5'	2.14	0.46
1:CA:1515:G:O2'	1:CA:1516:G:H5'	2.15	0.46
1:CA:465:A:O2'	1:CA:466:A:H2'	2.16	0.46
1:CA:845:A:H3'	1:CA:846:G:H8	1.80	0.46
1:CA:864:A:H2'	1:CA:865:A:C8	2.50	0.46
1:CA:13:U:C1'	1:CA:914:A:H5''	2.39	0.46
18:CB:172:ILE:H	18:CB:172:ILE:HG13	1.47	0.46
18:CB:223:GLY:C	18:CB:225:SER:H	2.19	0.46
18:CB:27:LYS:C	18:CB:29:PHE:H	2.19	0.46
2:CC:13:ILE:O	2:CC:15:LYS:N	2.49	0.46
2:CC:116:ALA:HB1	2:CC:197:VAL:HG11	1.98	0.46
2:CC:6:PRO:HG2	2:CC:183:TYR:CD2	2.50	0.46
3:CD:154:VAL:HG23	3:CD:155:LYS:N	2.31	0.46
6:CG:134:VAL:CG1	6:CG:137:ARG:HH21	2.27	0.46
10:CK:31:VAL:O	10:CK:43:TRP:HA	2.16	0.46
10:CK:95:THR:HG23	10:CK:96:ILE:H	1.80	0.46
12:CM:2:ARG:H	12:CM:2:ARG:CD	2.28	0.46
14:CQ:17:GLU:O	14:CQ:18:LYS:HB2	2.15	0.46
15:CR:41:SER:HB3	15:CR:51:GLN:CD	2.35	0.46
15:CR:32:ILE:HD11	15:CR:58:ILE:HG23	1.97	0.46
33:D1:50:GLU:O	33:D1:51:ALA:HB2	2.15	0.46
23:DB:1082:U:C2	23:DB:1086:A:N1	2.83	0.46
23:DB:1250:G:OP2	37:DL:21:ARG:NH2	2.48	0.46
23:DB:1729:U:H5	23:DB:1731:G:H22	1.63	0.46
23:DB:1799:G:N2	23:DB:1818:U:O2'	2.49	0.46
23:DB:235:U:H2'	23:DB:236:C:H6	1.81	0.46
23:DB:2511:U:H2'	23:DB:2512:C:C6	2.51	0.46
23:DB:2570:G:H2'	23:DB:2571:U:O4'	2.16	0.46
23:DB:2654:A:H61	23:DB:2665:A:H3'	1.80	0.46
23:DB:638:G:H2'	23:DB:639:U:C6	2.51	0.46
23:DB:685:A:H1'	23:DB:688:U:O4	2.15	0.46
25:DC:171:VAL:HG23	25:DC:185:ALA:CB	2.46	0.46
25:DC:202:ARG:HH11	25:DC:213:ARG:NE	2.14	0.46
40:DH:68:ARG:HA	40:DH:68:ARG:NE	2.31	0.46
28:DP:30:TRP:HD1	28:DP:39:LEU:HG	1.79	0.46
44:DQ:56:PHE:O	44:DQ:59:LEU:HB3	2.15	0.46
49:DR:2:TYR:CD1	49:DR:42:ALA:HB2	2.50	0.46
23:DB:96:C:H4'	39:DX:41:HIS:CG	2.51	0.46
30:DY:12:ALA:HB2	30:DY:53:MET:SD	2.55	0.46
30:DY:7:THR:HA	30:DY:34:THR:HA	1.98	0.46
51:DZ:70:GLU:HG3	51:DZ:71:LEU:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1012:A:H2'	1:AA:1013:G:O4'	2.16	0.46
1:AA:1050:G:C2'	1:AA:1051:C:H5'	2.45	0.46
1:AA:121:U:H3'	1:AA:121:U:OP1	2.16	0.46
1:AA:317:U:H2'	1:AA:318:G:H8	1.81	0.46
1:AA:692:U:H2'	1:AA:694:A:OP2	2.16	0.46
18:AB:130:LYS:N	18:AB:130:LYS:HD2	2.30	0.46
2:AC:151:GLU:HB3	2:AC:198:LYS:HB2	1.98	0.46
3:AD:59:LYS:HG2	3:AD:63:ILE:HD11	1.97	0.46
5:AF:53:LYS:NZ	5:AF:53:LYS:H	2.14	0.46
6:AG:22:LEU:HD23	6:AG:22:LEU:O	2.15	0.46
6:AG:92:PRO:HG2	6:AG:93:VAL:H	1.80	0.46
7:AH:99:GLY:CA	7:AH:129:ALA:HA	2.45	0.46
8:AI:35:GLU:CD	8:AI:39:GLY:HA2	2.36	0.46
10:AK:31:VAL:O	10:AK:43:TRP:HA	2.16	0.46
10:AK:69:CYS:SG	10:AK:70:ALA:N	2.89	0.46
11:AL:49:ARG:CB	11:AL:89:LEU:HD11	2.44	0.46
1:AA:947:G:H4'	12:AM:107:THR:OG1	2.16	0.46
12:AM:89:ARG:HH11	12:AM:94:LEU:HB3	1.80	0.46
21:AN:9:GLU:OE1	21:AN:59:GLN:HB3	2.15	0.46
14:AQ:20:ILE:HG12	14:AQ:52:CYS:SG	2.56	0.46
22:BA:37:C:H3'	22:BA:38:C:H6	1.81	0.46
23:BB:1328:A:H2'	23:BB:1330:C:C4	2.51	0.46
23:BB:159:G:O2'	23:BB:160:A:H5''	2.16	0.46
23:BB:1739:A:H2'	23:BB:1740:G:C8	2.50	0.46
23:BB:1793:C:H2'	23:BB:1794:A:C8	2.51	0.46
23:BB:185:G:H2'	23:BB:186:G:O4'	2.16	0.46
23:BB:1977:A:H2'	23:BB:1978:A:C8	2.51	0.46
23:BB:2027:G:O2'	23:BB:2028:U:H5'	2.16	0.46
23:BB:2570:G:H2'	23:BB:2571:U:O4'	2.15	0.46
23:BB:274:C:H2'	23:BB:275:C:C6	2.51	0.46
23:BB:359:G:H2'	23:BB:360:U:H5'	1.96	0.46
23:BB:437:U:H2'	23:BB:438:G:C8	2.50	0.46
23:BB:592:A:H2'	23:BB:593:U:C6	2.51	0.46
23:BB:902:C:H2'	23:BB:903:C:C6	2.51	0.46
23:BB:962:G:H2'	23:BB:963:U:C6	2.50	0.46
26:BD:183:GLU:N	26:BD:183:GLU:OE1	2.49	0.46
26:BD:62:LYS:HD2	26:BD:62:LYS:N	2.30	0.46
29:BE:1:MET:O	29:BE:13:THR:HA	2.15	0.46
47:BF:155:ILE:HG22	47:BF:156:THR:N	2.30	0.46
23:BB:2309:A:H61	47:BF:75:GLY:CA	2.28	0.46
48:BG:155:PRO:HA	48:BG:170:THR:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:5:LEU:HG	40:BH:17:ASP:O	2.16	0.46
41:BJ:20:ALA:HB2	41:BJ:28:LEU:CD1	2.45	0.46
41:BJ:34:ARG:HG3	41:BJ:39:LYS:HB3	1.96	0.46
37:BL:124:GLY:N	37:BL:143:GLU:CG	2.75	0.46
37:BL:98:ALA:O	37:BL:99:ASN:C	2.53	0.46
38:BM:120:ALA:O	38:BM:123:LYS:HG2	2.16	0.46
38:BM:26:VAL:HB	38:BM:104:GLU:OE2	2.16	0.46
43:BO:47:VAL:HG12	43:BO:48:LEU:N	2.26	0.46
43:BO:75:GLY:O	43:BO:78:VAL:HG23	2.16	0.46
28:BP:95:LYS:HG3	28:BP:97:TYR:HE1	1.80	0.46
44:BQ:7:VAL:HG23	44:BQ:8:ILE:N	2.30	0.46
44:BQ:64:ILE:HD12	44:BQ:95:ALA:CB	2.46	0.46
45:BS:30:SER:HA	45:BS:33:LEU:HD12	1.97	0.46
35:BV:11:GLU:CD	35:BV:16:ALA:HB1	2.35	0.46
35:BV:83:LYS:O	35:BV:85:LYS:N	2.49	0.46
52:BW:24:ARG:CD	52:BW:65:LYS:HG2	2.46	0.46
52:BW:9:THR:OG1	52:BW:10:ARG:N	2.46	0.46
51:BZ:27:ARG:O	51:BZ:28:ARG:CB	2.64	0.46
51:BZ:27:ARG:HG3	51:BZ:28:ARG:H	1.80	0.46
1:CA:1039:G:O2'	1:CA:1040:U:H5'	2.16	0.46
1:CA:20:U:H2'	1:CA:21:G:O4'	2.15	0.46
1:CA:456:A:H2'	1:CA:457:G:H5''	1.98	0.46
1:CA:679:C:O2'	1:CA:680:C:H5'	2.15	0.46
1:CA:897:C:O2'	1:CA:898:G:H5'	2.16	0.46
1:CA:244:U:O4	1:CA:906:A:H1'	2.15	0.46
1:CA:95:C:O2	1:CA:95:C:H2'	2.14	0.46
2:CC:112:ALA:N	2:CC:201:ILE:HD11	2.31	0.46
2:CC:138:GLN:O	2:CC:142:ARG:HB3	2.16	0.46
2:CC:19:SER:N	2:CC:55:VAL:HG13	2.31	0.46
2:CC:56:ILE:O	2:CC:58:ARG:HG3	2.16	0.46
3:CD:160:LEU:CD1	3:CD:160:LEU:H	2.24	0.46
3:CD:59:LYS:HG2	3:CD:63:ILE:HD11	1.98	0.46
4:CE:56:PRO:HA	4:CE:59:ILE:HD12	1.98	0.46
5:CF:53:LYS:C	5:CF:54:LEU:HD22	2.36	0.46
11:CL:85:ARG:CB	11:CL:93:ARG:HA	2.46	0.46
12:CM:108:ARG:HG3	12:CM:108:ARG:NH1	2.30	0.46
16:CS:64:GLU:H	16:CS:64:GLU:CD	2.17	0.46
23:DB:1197:G:O2'	23:DB:1198:U:H5'	2.15	0.46
23:DB:1316:U:H2'	23:DB:1317:G:H8	1.81	0.46
23:DB:1338:G:H4'	50:DT:18:GLU:HG3	1.97	0.46
23:DB:1429:G:H2'	23:DB:1430:G:C8	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:150:U:O2'	23:DB:151:C:H5'	2.15	0.46
23:DB:2800:A:H8	23:DB:2800:A:O5'	1.99	0.46
23:DB:574:A:H4'	23:DB:575:A:H5''	1.97	0.46
23:DB:672:C:H2'	23:DB:673:C:C6	2.51	0.46
23:DB:675:A:N3	23:DB:2443:C:O2'	2.49	0.46
26:DD:5:VAL:HG11	26:DD:80:TRP:CE3	2.51	0.46
29:DE:58:LYS:O	29:DE:60:TRP:N	2.49	0.46
47:DF:45:ASP:O	47:DF:47:LYS:N	2.43	0.46
47:DF:52:ALA:O	47:DF:56:LEU:HB2	2.14	0.46
48:DG:17:LYS:HE3	48:DG:18:ILE:N	2.31	0.46
24:DI:23:VAL:HG12	24:DI:24:GLY:N	2.31	0.46
23:DB:636:G:C6	37:DL:111:ILE:HD12	2.51	0.46
37:DL:80:SER:N	37:DL:113:ALA:HB3	2.30	0.46
42:DN:47:VAL:O	42:DN:51:LEU:HD22	2.15	0.46
43:DO:50:ALA:O	43:DO:51:ALA:HB2	2.15	0.46
28:DP:29:VAL:HA	28:DP:79:VAL:O	2.16	0.46
28:DP:63:ILE:HG12	28:DP:68:GLY:CA	2.46	0.46
45:DS:45:VAL:O	45:DS:48:LYS:HB3	2.16	0.46
45:DS:66:ILE:CA	45:DS:69:LEU:HD13	2.44	0.46
50:DT:53:VAL:HG11	50:DT:87:LEU:HD22	1.98	0.46
46:DU:10:VAL:HG22	46:DU:35:VAL:HG11	1.98	0.46
35:DV:56:PHE:O	35:DV:61:LEU:HD21	2.15	0.46
23:DB:2331:G:C4'	52:DW:39:GLN:HA	2.46	0.46
30:DY:35:VAL:HG22	30:DY:36:GLU:N	2.31	0.46
1:AA:1077:G:N2	1:AA:1080:A:OP2	2.48	0.46
1:AA:1333:A:C2	1:AA:1334:G:H1'	2.50	0.46
3:AD:155:LYS:HA	3:AD:158:LEU:HD12	1.98	0.46
4:AE:113:VAL:CG2	4:AE:114:LEU:H	2.20	0.46
4:AE:82:HIS:HB2	4:AE:83:PRO:HD2	1.97	0.46
5:AF:74:LEU:HG	5:AF:78:PHE:CE1	2.51	0.46
6:AG:13:PRO:HA	6:AG:23:ALA:HB2	1.97	0.46
8:AI:20:ILE:HA	8:AI:62:LEU:CD1	2.41	0.46
10:AK:14:GLN:HA	10:AK:76:TYR:O	2.16	0.46
20:AO:57:LEU:HD23	23:BB:715:A:C8	2.50	0.46
12:AM:78:ARG:NH1	16:AS:64:GLU:HB2	2.31	0.46
31:B0:18:HIS:C	31:B0:20:ALA:H	2.19	0.46
33:B1:18:HIS:CE1	33:B1:40:PRO:HD2	2.51	0.46
23:BB:1351:C:O2'	23:BB:1571:A:H1'	2.15	0.46
23:BB:1504:A:H2'	23:BB:1505:A:C8	2.51	0.46
23:BB:1590:A:H2'	23:BB:1591:A:C8	2.50	0.46
23:BB:1716:U:H2'	23:BB:1717:A:H8	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1915:U:H3'	23:BB:1916:A:H8	1.81	0.46
23:BB:2745:C:H2'	23:BB:2746:U:C6	2.50	0.46
23:BB:2863:C:O2'	23:BB:2864:G:H5'	2.16	0.46
23:BB:2884:U:C6	31:B0:49:ARG:HG2	2.51	0.46
23:BB:338:G:N2	23:BB:339:U:H1'	2.31	0.46
23:BB:969:G:H2'	23:BB:970:U:O4'	2.16	0.46
25:BC:116:GLN:O	25:BC:117:SER:HB3	2.15	0.46
25:BC:76:VAL:HA	25:BC:114:GLN:HA	1.97	0.46
26:BD:92:VAL:O	26:BD:94:GLN:N	2.49	0.46
47:BF:7:TYR:O	47:BF:11:VAL:HB	2.16	0.46
47:BF:37:MET:HA	47:BF:151:LEU:HA	1.97	0.46
47:BF:49:LEU:HD21	47:BF:84:ILE:O	2.15	0.46
47:BF:4:HIS:O	47:BF:7:TYR:HB3	2.15	0.46
48:BG:17:LYS:HB3	48:BG:24:THR:O	2.16	0.46
48:BG:7:PRO:O	48:BG:8:VAL:HB	2.15	0.46
24:BI:135:MET:HG3	24:BI:137:LEU:HG	1.98	0.46
41:BJ:115:GLY:HA2	41:BJ:118:MET:HE3	1.98	0.46
38:BM:114:ARG:HG3	38:BM:130:PHE:CD1	2.51	0.46
28:BP:32:VAL:HG12	28:BP:33:GLU:O	2.16	0.46
28:BP:4:ILE:C	28:BP:6:GLN:H	2.18	0.46
44:BQ:67:ALA:HB1	44:BQ:105:PHE:CZ	2.51	0.46
49:BR:2:TYR:HB2	49:BR:42:ALA:N	2.31	0.46
45:BS:57:ASN:HD22	45:BS:57:ASN:N	2.14	0.46
50:BT:51:PHE:HB3	50:BT:53:VAL:HG23	1.97	0.46
46:BU:11:ILE:HG21	46:BU:78:LYS:CE	2.46	0.46
35:BV:16:ALA:HA	35:BV:19:ARG:CZ	2.45	0.46
52:BW:9:THR:HG23	52:BW:10:ARG:CD	2.45	0.46
1:CA:1173:U:H2'	1:CA:1174:G:O4'	2.15	0.46
1:CA:18:C:H1'	1:CA:1079:G:N2	2.28	0.46
1:CA:425:G:H2'	1:CA:426:U:C6	2.51	0.46
1:CA:509:A:C6	1:CA:510:A:N1	2.84	0.46
1:CA:662:U:O2'	1:CA:836:G:H5''	2.16	0.46
2:CC:31:ASN:ND2	2:CC:58:ARG:HE	2.12	0.46
2:CC:68:HIS:O	2:CC:105:VAL:HG21	2.16	0.46
1:CA:8:A:H1'	4:CE:107:GLY:HA2	1.97	0.46
4:CE:33:THR:HB	4:CE:49:TYR:OH	2.16	0.46
9:CJ:81:GLU:N	9:CJ:82:LYS:NZ	2.64	0.46
10:CK:85:VAL:O	10:CK:111:ASP:HA	2.15	0.46
16:CS:30:LEU:HD12	16:CS:48:ILE:HG12	1.96	0.46
23:DB:2348:U:H5'	33:D1:20:TYR:OH	2.15	0.46
34:D3:49:VAL:HG11	34:D3:54:LEU:CD1	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1275:A:H2'	23:DB:1276:A:O4'	2.16	0.46
23:DB:1410:G:H2'	23:DB:1411:U:C6	2.51	0.46
23:DB:1590:A:H2'	23:DB:1591:A:C8	2.50	0.46
23:DB:1797:G:O2'	23:DB:1798:U:H5'	2.16	0.46
23:DB:2064:C:H2'	23:DB:2065:C:H6	1.76	0.46
23:DB:2267:A:C3'	23:DB:2267:A:H8	2.25	0.46
23:DB:278:A:H5''	23:DB:279:A:OP1	2.16	0.46
23:DB:510:C:H2'	23:DB:511:U:O4'	2.16	0.46
23:DB:615:U:O2	29:DE:35:TYR:HA	2.15	0.46
25:DC:76:VAL:HG12	25:DC:114:GLN:HB2	1.97	0.46
25:DC:16:VAL:CB	25:DC:203:VAL:HG11	2.45	0.46
47:DF:7:TYR:O	47:DF:11:VAL:HB	2.16	0.46
23:DB:1099:G:C8	24:DI:3:LYS:CB	2.99	0.46
27:DK:61:VAL:HG11	27:DK:112:PHE:CZ	2.50	0.46
42:DN:62:ASN:N	42:DN:62:ASN:ND2	2.63	0.46
44:DQ:48:ASP:C	44:DQ:50:ARG:N	2.69	0.46
30:DY:43:ILE:HG13	30:DY:44:ARG:N	2.30	0.46
51:DZ:32:ASN:HB3	51:DZ:34:HIS:NE2	2.31	0.46
1:AA:1245:C:O2'	1:AA:1246:A:H5'	2.16	0.46
1:AA:1317:C:C2	1:AA:1318:A:H1'	2.51	0.46
1:AA:415:A:OP1	1:AA:415:A:H4'	2.16	0.46
1:AA:520:A:N1	1:AA:536:C:H1'	2.31	0.46
1:AA:292:G:O2'	1:AA:609:A:N6	2.49	0.46
1:AA:656:G:H2'	1:AA:657:U:H6	1.81	0.46
1:AA:796:C:O3'	10:AK:126:ARG:NH2	2.49	0.46
18:AB:117:GLU:HG3	18:AB:140:LEU:HD22	1.97	0.46
3:AD:160:LEU:H	3:AD:160:LEU:CD1	2.23	0.46
6:AG:87:PRO:CG	6:AG:144:ALA:HA	2.46	0.46
7:AH:14:ARG:HG3	7:AH:15:ASN:N	2.31	0.46
10:AK:23:HIS:HB3	10:AK:30:ILE:CG1	2.42	0.46
11:AL:6:LEU:HB2	14:AQ:33:TYR:CE1	2.51	0.46
20:AO:61:SER:O	20:AO:64:ARG:HB3	2.15	0.46
33:B1:38:PHE:HB2	33:B1:45:HIS:CE1	2.51	0.46
23:BB:1416:G:HO2'	23:BB:1417:C:H6	1.63	0.46
23:BB:1435:G:H2'	23:BB:1436:G:H8	1.80	0.46
23:BB:1688:U:O2	23:BB:1700:A:H5'	2.15	0.46
23:BB:1790:C:H2'	23:BB:1791:A:C8	2.51	0.46
23:BB:2348:U:H5'	33:B1:20:TYR:HH	1.81	0.46
23:BB:2722:G:H2'	23:BB:2723:C:H6	1.79	0.46
23:BB:2815:C:H2'	23:BB:2816:G:H8	1.80	0.46
23:BB:428:A:O2'	23:BB:429:A:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:549:G:H5''	23:BB:550:C:C5	2.51	0.46
25:BC:62:ARG:HB3	25:BC:63:ILE:H	1.43	0.46
26:BD:133:THR:HG23	26:BD:134:HIS:CD2	2.51	0.46
23:BB:2050:C:H1'	26:BD:161:MET:CE	2.46	0.46
29:BE:170:ARG:NH2	29:BE:176:ASP:HB2	2.30	0.46
47:BF:31:GLU:HG3	47:BF:32:LYS:HG2	1.97	0.46
48:BG:6:ALA:HB3	48:BG:68:ARG:CG	2.40	0.46
40:BH:117:LEU:CG	40:BH:130:VAL:HG13	2.45	0.46
40:BH:46:PHE:O	40:BH:47:PHE:HD2	1.99	0.46
40:BH:94:ILE:HB	40:BH:128:HIS:HE2	1.79	0.46
42:BN:103:ARG:HB2	42:BN:110:MET:CE	2.46	0.46
50:BT:39:THR:C	50:BT:41:ALA:H	2.20	0.46
35:BV:1:MET:O	35:BV:2:PHE:HB2	2.16	0.46
1:CA:1071:C:C2	1:CA:1072:G:C8	3.04	0.46
1:CA:1390:U:H2'	1:CA:1391:U:C6	2.51	0.46
1:CA:410:G:H1'	1:CA:432:A:H61	1.81	0.46
1:CA:628:G:O2'	1:CA:629:A:H5'	2.15	0.46
1:CA:62:U:H2'	1:CA:63:C:C6	2.51	0.46
1:CA:682:G:O2'	1:CA:683:G:H5'	2.16	0.46
1:CA:949:A:O2'	1:CA:950:U:H5'	2.16	0.46
1:CA:950:U:H2'	1:CA:951:G:C8	2.50	0.46
2:CC:152:VAL:H	2:CC:165:GLU:HB3	1.81	0.46
1:CA:1078:U:H4'	4:CE:137:ARG:CZ	2.46	0.46
7:CH:34:ALA:HB1	7:CH:109:VAL:HB	1.98	0.46
1:CA:779:C:H5''	10:CK:123:PRO:HB3	1.97	0.46
20:CO:49:ASP:O	20:CO:50:HIS:C	2.53	0.46
20:CO:85:LEU:HB3	20:CO:87:LEU:HG	1.97	0.46
1:CA:636:U:OP1	14:CQ:5:ARG:HD2	2.15	0.46
16:CS:10:ILE:O	16:CS:11:ASP:C	2.54	0.46
16:CS:64:GLU:O	16:CS:66:VAL:HG22	2.16	0.46
33:D1:32:LYS:NZ	33:D1:52:LYS:HG2	2.31	0.46
32:D4:13:ASN:HB2	32:D4:27:CYS:SG	2.56	0.46
23:DB:1098:A:C4	24:DI:3:LYS:O	2.68	0.46
23:DB:143:C:H2'	23:DB:144:A:C8	2.51	0.46
23:DB:1710:G:O2'	23:DB:1711:A:H5'	2.15	0.46
23:DB:1818:U:HO2'	23:DB:1819:A:P	2.38	0.46
23:DB:1859:U:H6	23:DB:1859:U:O5'	1.99	0.46
23:DB:2181:U:H2'	23:DB:2182:U:C6	2.51	0.46
23:DB:2353:G:N3	52:DW:30:VAL:HG13	2.31	0.46
23:DB:2418:A:H2'	23:DB:2419:U:O4'	2.16	0.46
23:DB:2588:G:H3'	56:DB:3587:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2599:G:H2'	23:DB:2600:A:C8	2.51	0.46
23:DB:2723:C:H5''	42:DN:1:MET:CE	2.46	0.46
23:DB:49:A:OP1	23:DB:51:G:H5'	2.16	0.46
23:DB:553:G:C2'	23:DB:554:U:H5'	2.46	0.46
23:DB:708:G:H2'	23:DB:709:U:C6	2.50	0.46
23:DB:95:A:H4'	39:DX:38:GLN:O	2.14	0.46
25:DC:68:ARG:HB3	25:DC:128:THR:OG1	2.16	0.46
25:DC:196:ASN:HD21	25:DC:199:HIS:HB2	1.78	0.46
25:DC:41:GLY:HA3	25:DC:53:ILE:CG2	2.37	0.46
26:DD:114:LYS:HD2	26:DD:116:LYS:CE	2.45	0.46
26:DD:46:ARG:HG3	26:DD:46:ARG:HH11	1.80	0.46
29:DE:97:ASN:ND2	29:DE:100:MET:HG3	2.30	0.46
47:DF:116:LEU:O	47:DF:177:ARG:HB2	2.15	0.46
47:DF:134:GLN:C	47:DF:136:ILE:N	2.69	0.46
38:DM:113:ALA:O	38:DM:116:ALA:HB3	2.16	0.46
42:DN:2:ARG:HB3	42:DN:2:ARG:CZ	2.45	0.46
49:DR:14:VAL:HG21	49:DR:98:ILE:HD13	1.98	0.46
50:DT:10:VAL:O	50:DT:12:ARG:N	2.48	0.46
50:DT:55:VAL:CA	50:DT:87:LEU:HA	2.30	0.46
1:AA:1007:U:H2'	1:AA:1008:U:H6	1.79	0.46
1:AA:1102:A:O2'	1:AA:1103:C:H5'	2.16	0.46
1:AA:1308:U:OP1	12:AM:96:VAL:HG23	2.15	0.46
1:AA:56:U:H2'	1:AA:57:G:H8	1.81	0.46
1:AA:864:A:H2'	1:AA:865:A:C8	2.51	0.46
1:AA:986:U:H1'	16:AS:54:ARG:HB3	1.97	0.46
3:AD:155:LYS:HG2	3:AD:156:ALA:H	1.81	0.46
4:AE:129:SER:HA	56:AE:201:HOH:O	2.16	0.46
4:AE:150:GLU:H	4:AE:150:GLU:CD	2.19	0.46
6:AG:136:LYS:O	6:AG:140:VAL:HG23	2.16	0.46
8:AI:53:LEU:O	8:AI:53:LEU:HD13	2.15	0.46
8:AI:66:VAL:O	8:AI:67:LYS:HG3	2.16	0.46
10:AK:121:ARG:HH12	19:AU:37:TYR:HE2	1.64	0.46
10:AK:27:ASN:O	10:AK:56:LYS:HE2	2.16	0.46
11:AL:23:LEU:HD22	11:AL:58:ASN:HB2	1.98	0.46
11:AL:35:ARG:HA	11:AL:35:ARG:CZ	2.46	0.46
11:AL:5:GLN:HA	11:AL:8:ARG:NH2	2.31	0.46
21:AN:25:GLU:O	21:AN:29:ILE:HG13	2.16	0.46
23:BB:2056:G:H21	31:B0:1:ALA:HA	1.82	0.46
31:B0:41:HIS:HB3	42:BN:99:LYS:HB2	1.97	0.46
36:B2:13:ASN:N	36:B2:13:ASN:HD22	2.14	0.46
22:BA:111:U:H2'	22:BA:112:G:H8	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1119:U:H2'	23:BB:1120:G:H8	1.81	0.46
23:BB:1571:A:H2'	23:BB:1572:A:H8	1.81	0.46
23:BB:1863:G:H2'	23:BB:1864:U:O4'	2.16	0.46
23:BB:2144:G:N2	23:BB:2146:C:H5'	2.31	0.46
23:BB:2201:G:H2'	23:BB:2202:U:H6	1.80	0.46
23:BB:2305:U:H2'	23:BB:2306:C:O4'	2.16	0.46
23:BB:282:A:O2'	23:BB:283:G:H5'	2.16	0.46
23:BB:43:G:H2'	23:BB:44:A:O4'	2.16	0.46
23:BB:690:G:H2'	23:BB:691:C:O4'	2.16	0.46
26:BD:159:LYS:NZ	26:BD:160:LYS:N	2.64	0.46
26:BD:168:GLU:HG3	26:BD:170:VAL:CG1	2.46	0.46
26:BD:34:VAL:HA	26:BD:50:VAL:HG12	1.98	0.46
29:BE:46:GLN:HG3	29:BE:87:ALA:HB3	1.98	0.46
47:BF:69:ALA:O	47:BF:70:ARG:HB2	2.16	0.46
48:BG:146:ASP:O	48:BG:150:TYR:HD1	1.99	0.46
40:BH:104:THR:HG23	40:BH:105:ALA:H	1.79	0.46
40:BH:67:ALA:HA	40:BH:138:VAL:HG11	1.98	0.46
41:BJ:64:VAL:HG22	41:BJ:68:LYS:HB2	1.97	0.46
37:BL:10:GLU:C	37:BL:12:SER:H	2.19	0.46
37:BL:61:LEU:N	37:BL:61:LEU:CD1	2.79	0.46
38:BM:71:LYS:O	38:BM:92:TRP:HA	2.16	0.46
28:BP:30:TRP:HD1	28:BP:39:LEU:HG	1.81	0.46
26:BD:15:PHE:CE2	28:BP:77:SER:HB2	2.50	0.46
50:BT:40:LYS:C	50:BT:43:ILE:HG22	2.36	0.46
46:BU:24:VAL:HA	46:BU:35:VAL:HA	1.97	0.46
39:BX:45:GLN:O	39:BX:47:ARG:N	2.45	0.46
39:BX:44:LYS:HD2	39:BX:48:ARG:HH12	1.80	0.46
30:BY:7:THR:HA	30:BY:34:THR:HA	1.98	0.46
1:CA:1403:C:O2'	1:CA:1404:C:H5'	2.16	0.46
1:CA:1495:U:O2'	1:CA:1496:C:H5'	2.16	0.46
1:CA:218:U:H2'	1:CA:219:U:O4'	2.16	0.46
1:CA:286:C:H2'	1:CA:287:U:C6	2.51	0.46
1:CA:306:A:H2'	1:CA:307:C:H5'	1.98	0.46
1:CA:502:A:H2'	1:CA:503:C:H6	1.81	0.46
1:CA:763:G:H2'	1:CA:764:C:H6	1.81	0.46
18:CB:114:LYS:HZ3	18:CB:151:LYS:CD	2.29	0.46
4:CE:54:GLU:HG2	4:CE:56:PRO:HD2	1.98	0.46
4:CE:56:PRO:HA	4:CE:59:ILE:CD1	2.46	0.46
6:CG:91:ARG:HD2	6:CG:91:ARG:N	2.31	0.46
8:CI:122:ARG:HH11	8:CI:122:ARG:HG3	1.80	0.46
8:CI:64:ILE:HG21	8:CI:78:ILE:HG12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CL:23:LEU:HD22	11:CL:58:ASN:HB2	1.98	0.46
20:CO:70:LEU:CD1	20:CO:78:TYR:H	2.29	0.46
11:CL:6:LEU:HB2	14:CQ:33:TYR:CE1	2.50	0.46
17:CT:25:SER:O	17:CT:29:THR:HG23	2.16	0.46
23:DB:1092:C:H2'	23:DB:1093:G:C5'	2.45	0.46
23:DB:1174:U:O3'	23:DB:1175:A:H8	1.99	0.46
23:DB:1219:U:H2'	23:DB:1220:G:H8	1.80	0.46
23:DB:1241:A:O4'	23:DB:1241:A:N3	2.49	0.46
23:DB:1295:C:H2'	23:DB:1296:G:C8	2.51	0.46
23:DB:1383:A:H2	23:DB:1405:U:O2	1.99	0.46
23:DB:1439:A:N3	23:DB:1553:A:C6	2.84	0.46
23:DB:1469:A:H2'	23:DB:1470:A:H8	1.81	0.46
23:DB:1706:C:H2'	23:DB:1757:A:OP2	2.16	0.46
23:DB:1751:U:H2'	23:DB:1752:C:C6	2.51	0.46
23:DB:2361:G:H2'	23:DB:2362:C:H6	1.80	0.46
23:DB:2443:C:O2'	23:DB:2444:G:H5'	2.16	0.46
23:DB:245:G:O2'	23:DB:246:C:H5'	2.16	0.46
23:DB:2527:C:O2'	23:DB:2528:U:H5'	2.16	0.46
23:DB:2636:C:H2'	23:DB:2637:U:H6	1.81	0.46
23:DB:1750:G:O2'	23:DB:2860:A:H2	1.99	0.46
23:DB:2888:C:H2'	23:DB:2889:C:H6	1.81	0.46
23:DB:107:G:H1'	23:DB:294:A:H1'	1.98	0.46
23:DB:311:A:H2	23:DB:331:C:H5'	1.79	0.46
23:DB:349:U:O2	23:DB:349:U:H2'	2.15	0.46
23:DB:57:C:H2'	23:DB:58:G:C8	2.50	0.46
23:DB:876:C:H3'	23:DB:877:A:C4'	2.46	0.46
25:DC:202:ARG:HB3	25:DC:203:VAL:H	1.67	0.46
26:DD:183:GLU:N	26:DD:183:GLU:CD	2.69	0.46
26:DD:187:LEU:O	26:DD:188:LEU:HD23	2.16	0.46
29:DE:17:THR:C	29:DE:19:PHE:H	2.20	0.46
41:DJ:131:ASN:O	41:DJ:132:HIS:ND1	2.49	0.46
28:DP:19:PHE:HB2	28:DP:50:ARG:NH1	2.30	0.46
44:DQ:64:ILE:HD12	44:DQ:95:ALA:CB	2.45	0.46
49:DR:58:VAL:HG22	49:DR:59:ILE:N	2.22	0.46
46:DU:3:LYS:HD3	46:DU:82:VAL:HG21	1.98	0.46
35:DV:83:LYS:O	35:DV:85:LYS:N	2.49	0.46
1:AA:1114:C:O2'	1:AA:1115:U:H5'	2.16	0.45
1:AA:1217:C:H2'	1:AA:1218:C:H6	1.81	0.45
1:AA:1244:G:H2'	1:AA:1245:C:O4'	2.16	0.45
1:AA:1299:A:H2'	1:AA:1301:U:C1'	2.47	0.45
1:AA:386:C:O2'	1:AA:387:U:H5'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:409:U:OP1	3:AD:21:LYS:HG3	2.15	0.45
1:AA:429:U:C1'	1:AA:430:A:H5''	2.45	0.45
1:AA:778:G:O2'	1:AA:779:C:H5'	2.16	0.45
1:AA:956:U:C2'	1:AA:957:U:H5'	2.47	0.45
2:AC:152:VAL:HG13	2:AC:197:VAL:HG22	1.98	0.45
2:AC:61:LYS:O	2:AC:96:VAL:HB	2.16	0.45
2:AC:78:LYS:C	2:AC:80:GLY:H	2.20	0.45
3:AD:146:GLU:CD	3:AD:146:GLU:H	2.19	0.45
4:AE:91:SER:OG	4:AE:135:VAL:HG22	2.15	0.45
6:AG:144:ALA:O	6:AG:145:GLU:HB3	2.16	0.45
8:AI:64:ILE:HG22	8:AI:66:VAL:H	1.81	0.45
12:AM:22:TYR:HB2	12:AM:65:GLU:HA	1.98	0.45
21:AN:61:ASN:HB3	21:AN:72:PHE:CD2	2.51	0.45
20:AO:70:LEU:HD12	20:AO:78:TYR:HB2	1.98	0.45
15:AR:41:SER:HB3	15:AR:51:GLN:HB2	1.98	0.45
19:AU:35:GLU:HB2	19:AU:37:TYR:CZ	2.51	0.45
34:B3:2:LYS:HG2	34:B3:2:LYS:H	1.45	0.45
23:BB:1006:C:H1'	41:BJ:108:MET:SD	2.57	0.45
23:BB:1389:G:O2'	23:BB:1390:U:H5'	2.16	0.45
23:BB:1468:U:H2'	23:BB:1522:A:N6	2.31	0.45
23:BB:1824:G:H2'	23:BB:1825:U:H6	1.82	0.45
23:BB:2269:G:H4'	52:BW:19:ARG:NH1	2.30	0.45
23:BB:2271:G:H2'	23:BB:2272:U:C6	2.51	0.45
23:BB:2297:A:N6	23:BB:2319:G:H1'	2.31	0.45
23:BB:2742:G:H2'	23:BB:2743:U:H6	1.80	0.45
23:BB:2740:A:N7	23:BB:2764:A:N7	2.64	0.45
23:BB:2805:C:O2'	23:BB:2806:C:H5'	2.16	0.45
23:BB:2840:C:H5''	42:BN:53:THR:CG2	2.46	0.45
23:BB:513:A:O5'	23:BB:513:A:H8	1.98	0.45
23:BB:527:C:O4'	23:BB:527:C:O2	2.30	0.45
25:BC:251:THR:O	25:BC:252:LYS:HD2	2.16	0.45
48:BG:112:VAL:O	48:BG:113:ASP:HB2	2.16	0.45
48:BG:108:PHE:HE1	48:BG:151:ARG:HH11	1.64	0.45
48:BG:90:GLY:HA3	48:BG:93:TYR:CZ	2.52	0.45
38:BM:53:MET:SD	38:BM:63:ILE:HG21	2.56	0.45
38:BM:19:GLY:CA	38:BM:98:PRO:HD2	2.37	0.45
22:BA:114:C:H1'	43:BO:47:VAL:HG21	1.98	0.45
28:BP:3:ILE:HG23	28:BP:4:ILE:N	2.31	0.45
28:BP:50:ARG:HB3	28:BP:57:ALA:N	2.31	0.45
27:BK:76:VAL:O	28:BP:71:ARG:HA	2.16	0.45
44:BQ:72:GLY:O	44:BQ:113:LYS:HE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BQ:115:ALA:C	44:BQ:117:ALA:H	2.20	0.45
44:BQ:31:TYR:O	44:BQ:32:ARG:C	2.54	0.45
44:BQ:93:ILE:O	44:BQ:96:ASP:HB3	2.15	0.45
44:BQ:9:ALA:O	44:BQ:11:ALA:N	2.49	0.45
50:BT:45:ALA:HA	50:BT:48:GLN:HG2	1.98	0.45
50:BT:69:ARG:HG3	50:BT:70:HIS:N	2.30	0.45
39:BX:39:GLN:HB3	39:BX:41:HIS:NE2	2.30	0.45
1:CA:1225:A:H5'	12:CM:101:THR:HG1	1.81	0.45
1:CA:1463:U:H2'	1:CA:1464:U:H6	1.81	0.45
1:CA:411:A:H62	1:CA:413:G:H21	1.64	0.45
1:CA:45:G:O2'	1:CA:46:G:H5'	2.16	0.45
1:CA:473:U:H2'	1:CA:474:G:H8	1.81	0.45
1:CA:492:C:H2'	1:CA:493:A:N3	2.31	0.45
1:CA:72:A:N6	1:CA:98:A:C2	2.83	0.45
1:CA:813:U:H5''	1:CA:816:A:N6	2.31	0.45
1:CA:88:U:H2'	1:CA:89:U:C6	2.51	0.45
2:CC:53:ARG:HB2	2:CC:68:HIS:CD2	2.52	0.45
3:CD:21:LYS:C	3:CD:23:GLY:H	2.20	0.45
3:CD:58:GLN:HA	3:CD:58:GLN:OE1	2.16	0.45
5:CF:38:ARG:HH21	5:CF:63:ASN:HD21	1.64	0.45
5:CF:55:HIS:HB2	5:CF:56:LYS:HE2	1.98	0.45
5:CF:77:THR:O	5:CF:81:ASN:HB2	2.16	0.45
6:CG:3:ARG:HB3	6:CG:4:ARG:H	1.53	0.45
6:CG:94:ARG:HD3	6:CG:98:LEU:HD12	1.98	0.45
4:CE:156:ARG:HB3	7:CH:43:GLY:O	2.16	0.45
8:CI:49:GLN:N	8:CI:50:PRO:HD2	2.31	0.45
8:CI:25:GLY:HA3	8:CI:57:VAL:CA	2.46	0.45
9:CJ:83:THR:HA	9:CJ:86:ALA:HB3	1.98	0.45
11:CL:79:ILE:HD12	11:CL:96:THR:HG22	1.97	0.45
21:CN:8:ARG:O	21:CN:12:ARG:HG3	2.16	0.45
22:DA:59:A:H2'	22:DA:60:C:H6	1.78	0.45
23:DB:1177:G:O5'	23:DB:1177:G:H8	1.99	0.45
23:DB:1198:U:H2'	23:DB:1199:U:H6	1.79	0.45
23:DB:1501:G:O2'	23:DB:1502:A:H5'	2.15	0.45
23:DB:1654:A:H61	23:DB:2049:G:P	2.39	0.45
23:DB:2246:G:H2'	23:DB:2247:A:C8	2.51	0.45
23:DB:2514:U:H2'	23:DB:2515:C:C6	2.51	0.45
23:DB:2634:A:H2'	23:DB:2635:A:H8	1.80	0.45
23:DB:2836:U:H2'	23:DB:2837:A:C8	2.51	0.45
23:DB:428:A:O2'	23:DB:429:A:H5'	2.16	0.45
23:DB:84:A:OP1	46:DU:2:ALA:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:849:A:H2'	23:DB:850:U:C6	2.51	0.45
23:DB:957:C:N4	23:DB:2459:A:C8	2.84	0.45
25:DC:143:VAL:HA	25:DC:189:ALA:CB	2.46	0.45
25:DC:62:ARG:HB3	25:DC:63:ILE:H	1.43	0.45
23:DB:1654:A:O2'	26:DD:118:PHE:CB	2.64	0.45
26:DD:36:GLN:O	26:DD:36:GLN:HG3	2.16	0.45
29:DE:128:ALA:O	29:DE:130:LYS:N	2.46	0.45
48:DG:155:PRO:CA	48:DG:170:THR:HA	2.46	0.45
40:DH:90:LEU:HD22	40:DH:122:LEU:O	2.16	0.45
40:DH:93:SER:C	40:DH:94:ILE:HD12	2.36	0.45
27:DK:115:ILE:HG23	27:DK:116:ILE:H	1.81	0.45
27:DK:62:VAL:HG12	27:DK:63:VAL:N	2.31	0.45
38:DM:31:PHE:HA	38:DM:131:VAL:O	2.16	0.45
49:DR:14:VAL:HG22	49:DR:15:SER:H	1.80	0.45
45:DS:24:ILE:HG12	45:DS:36:LEU:HD21	1.98	0.45
45:DS:49:LYS:C	45:DS:51:LEU:N	2.69	0.45
45:DS:50:VAL:HG12	45:DS:105:VAL:HG23	1.98	0.45
46:DU:35:VAL:HG23	46:DU:38:ILE:HG21	1.97	0.45
52:DW:43:LYS:O	52:DW:58:LEU:HD11	2.15	0.45
30:DY:6:ILE:HG22	30:DY:56:VAL:HA	1.98	0.45
1:AA:1111:A:H2	2:AC:176:THR:HG23	1.81	0.45
1:AA:55:A:OP2	1:AA:352:C:N4	2.49	0.45
1:AA:469:C:O2'	1:AA:470:C:H5'	2.17	0.45
1:AA:736:C:H2'	1:AA:737:C:H6	1.77	0.45
1:AA:738:C:H2'	1:AA:739:C:H6	1.81	0.45
1:AA:88:U:O2'	1:AA:89:U:C5	2.68	0.45
18:AB:185:ILE:O	18:AB:185:ILE:HG22	2.15	0.45
18:AB:78:ALA:HB1	18:AB:213:LEU:HD13	1.98	0.45
2:AC:100:ILE:O	2:AC:100:ILE:HG23	2.16	0.45
3:AD:117:VAL:HA	3:AD:122:ILE:HG13	1.96	0.45
4:AE:56:PRO:HA	4:AE:59:ILE:HD12	1.98	0.45
4:AE:61:LYS:HG3	4:AE:62:ALA:N	2.31	0.45
5:AF:53:LYS:C	5:AF:54:LEU:HD22	2.37	0.45
6:AG:74:VAL:HA	6:AG:86:VAL:C	2.36	0.45
7:AH:125:ILE:HG22	7:AH:126:CYS:SG	2.56	0.45
8:AI:113:LYS:HA	8:AI:120:ALA:HB2	1.97	0.45
8:AI:119:LYS:O	8:AI:121:ARG:N	2.49	0.45
8:AI:16:ALA:HB2	8:AI:66:VAL:HG23	1.98	0.45
21:AN:41:TRP:NE1	21:AN:43:ALA:HB3	2.32	0.45
21:AN:73:LEU:O	21:AN:77:GLY:N	2.49	0.45
21:AN:79:SER:OG	21:AN:82:LYS:HG2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AR:42:ARG:C	15:AR:44:THR:H	2.19	0.45
16:AS:17:LYS:HD2	16:AS:30:LEU:CD2	2.43	0.45
23:BB:1841:U:H2'	23:BB:1842:G:C8	2.51	0.45
23:BB:2221:G:O2'	23:BB:2222:C:H5'	2.16	0.45
23:BB:250:G:H2'	23:BB:251:A:C8	2.52	0.45
23:BB:2514:U:H2'	23:BB:2515:C:C6	2.52	0.45
23:BB:2700:A:H2'	23:BB:2701:U:H6	1.80	0.45
23:BB:494:G:O2'	23:BB:495:G:H5'	2.16	0.45
23:BB:710:U:H2'	23:BB:711:G:C8	2.51	0.45
23:BB:715:A:H2'	23:BB:716:A:O4'	2.16	0.45
23:BB:808:G:H2'	23:BB:809:G:O4'	2.16	0.45
23:BB:861:A:H2'	23:BB:862:G:O4'	2.16	0.45
26:BD:89:GLU:HG3	26:BD:94:GLN:OE1	2.16	0.45
48:BG:156:TYR:O	48:BG:157:LYS:HB2	2.16	0.45
40:BH:6:LEU:O	40:BH:15:LEU:HD22	2.15	0.45
40:BH:30:LEU:O	40:BH:35:LYS:HB2	2.15	0.45
24:BI:52:LEU:HD21	24:BI:81:LYS:NZ	2.31	0.45
24:BI:62:ALA:C	24:BI:64:ARG:H	2.20	0.45
37:BL:132:ARG:CA	37:BL:135:ILE:HG22	2.46	0.45
38:BM:32:GLY:HA3	38:BM:103:TYR:O	2.16	0.45
38:BM:41:LEU:O	38:BM:93:VAL:HA	2.16	0.45
42:BN:48:VAL:HG23	42:BN:49:GLU:N	2.32	0.45
43:BO:50:ALA:O	43:BO:51:ALA:HB2	2.17	0.45
49:BR:76:LYS:HB2	49:BR:85:LYS:HB2	1.99	0.45
45:BS:69:LEU:HD23	45:BS:107:VAL:HG21	1.97	0.45
50:BT:50:LEU:O	50:BT:51:PHE:HB2	2.15	0.45
35:BV:80:HIS:HA	35:BV:87:GLN:OE1	2.16	0.45
39:BX:22:LEU:HG	39:BX:23:ARG:HG2	1.98	0.45
51:BZ:18:ARG:HA	51:BZ:24:ALA:HA	1.98	0.45
1:CA:1390:U:H2'	1:CA:1391:U:H6	1.81	0.45
1:CA:764:C:O2'	1:CA:765:G:H5'	2.17	0.45
1:CA:85:U:HO2'	1:CA:86:G:P	2.39	0.45
18:CB:169:HIS:HD2	18:CB:170:ILE:H	1.63	0.45
18:CB:76:SER:OG	18:CB:77:GLU:N	2.49	0.45
2:CC:35:ASP:OD2	2:CC:56:ILE:HD12	2.16	0.45
8:CI:40:ARG:N	8:CI:44:ARG:HD3	2.32	0.45
8:CI:60:LEU:O	8:CI:60:LEU:HD12	2.16	0.45
10:CK:22:ILE:HD13	10:CK:95:THR:HG21	1.97	0.45
11:CL:24:GLU:C	11:CL:26:CYS:H	2.19	0.45
12:CM:49:GLU:HA	12:CM:52:ILE:CG2	2.45	0.45
20:CO:64:ARG:NH2	20:CO:88:ARG:HD2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:CQ:82:VAL:O	14:CQ:83:LEU:HD13	2.17	0.45
16:CS:49:ALA:HB1	16:CS:56:HIS:HB2	1.97	0.45
19:CU:20:ARG:H	19:CU:20:ARG:CD	2.30	0.45
36:D2:18:PHE:CE1	36:D2:22:MET:HG3	2.51	0.45
23:DB:1040:A:H2	23:DB:1115:G:H22	1.65	0.45
23:DB:1334:G:O2'	23:DB:1335:C:H5'	2.15	0.45
23:DB:1583:A:H4'	23:DB:1585:C:N3	2.31	0.45
23:DB:1688:U:O2	23:DB:1700:A:H5'	2.17	0.45
23:DB:1735:A:H2'	23:DB:1736:U:C6	2.51	0.45
23:DB:1863:G:H2'	23:DB:1864:U:O4'	2.15	0.45
23:DB:2271:G:H2'	23:DB:2272:U:C6	2.51	0.45
23:DB:2340:A:H2'	23:DB:2341:G:H8	1.81	0.45
23:DB:2708:G:O2'	23:DB:2709:G:H5'	2.17	0.45
23:DB:291:G:O2'	23:DB:292:U:H5'	2.16	0.45
23:DB:451:U:C2	23:DB:453:A:N7	2.84	0.45
23:DB:622:G:O2'	23:DB:623:C:H5'	2.16	0.45
25:DC:149:LYS:HG2	25:DC:152:GLN:OE1	2.16	0.45
26:DD:130:GLN:HE21	26:DD:130:GLN:HB3	1.56	0.45
26:DD:34:VAL:HG13	26:DD:94:GLN:H	1.77	0.45
29:DE:37:ALA:O	29:DE:39:ALA:N	2.43	0.45
47:DF:151:LEU:CD1	47:DF:153:ILE:HG13	2.46	0.45
47:DF:89:THR:O	47:DF:91:ARG:CZ	2.65	0.45
47:DF:89:THR:OG1	47:DF:91:ARG:NH2	2.49	0.45
48:DG:143:VAL:O	48:DG:147:LEU:HG	2.17	0.45
48:DG:84:LYS:HG3	48:DG:131:VAL:CB	2.44	0.45
40:DH:2:GLN:HB2	40:DH:39:ALA:HB2	1.97	0.45
24:DI:46:ASP:HA	24:DI:50:LYS:HE2	1.97	0.45
41:DJ:84:ILE:HG23	41:DJ:84:ILE:O	2.15	0.45
37:DL:98:ALA:O	37:DL:99:ASN:C	2.53	0.45
38:DM:17:ASN:HD21	38:DM:95:LEU:CG	2.29	0.45
38:DM:17:ASN:HD21	38:DM:95:LEU:HG	1.80	0.45
42:DN:70:THR:HB	42:DN:75:ILE:HD11	1.98	0.45
44:DQ:63:ARG:NH2	44:DQ:96:ASP:HA	2.32	0.45
49:DR:97:LYS:C	49:DR:98:ILE:HD12	2.37	0.45
50:DT:68:LYS:HD3	50:DT:68:LYS:N	2.31	0.45
1:AA:1236:A:H4'	1:AA:1304:G:H4'	1.98	0.45
1:AA:1270:G:H2'	1:AA:1271:A:C8	2.51	0.45
1:AA:1496:C:HO2'	1:AA:1517:G:H1	1.65	0.45
1:AA:332:G:H2'	1:AA:333:U:H6	1.81	0.45
1:AA:502:A:H2'	1:AA:503:C:H6	1.82	0.45
1:AA:533:A:H5''	56:AA:1892:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:791:G:C6	1:AA:792:A:N7	2.85	0.45
18:AB:8:MET:HB2	18:AB:11:ALA:HB3	1.97	0.45
2:AC:182:ASP:O	2:AC:201:ILE:HB	2.17	0.45
1:AA:8:A:N6	3:AD:53:GLN:HE22	2.08	0.45
3:AD:57:LYS:O	3:AD:61:ARG:HB2	2.17	0.45
11:AL:28:GLN:HB3	11:AL:28:GLN:HE21	1.53	0.45
21:AN:67:GLY:O	21:AN:69:PRO:HD3	2.15	0.45
13:AP:23:ASP:OD1	13:AP:25:ARG:HB2	2.17	0.45
13:AP:24:SER:C	13:AP:25:ARG:HD3	2.37	0.45
17:AT:66:ILE:HG23	17:AT:70:LYS:HD3	1.98	0.45
31:B0:55:ALA:O	31:B0:56:LYS:C	2.55	0.45
22:BA:29:A:H3'	22:BA:30:C:H6	1.81	0.45
22:BA:74:U:O2'	22:BA:75:G:H5'	2.16	0.45
23:BB:1035:U:H2'	23:BB:1036:G:C8	2.43	0.45
23:BB:1176:U:H2'	23:BB:1177:G:O4'	2.17	0.45
23:BB:1295:C:H2'	23:BB:1296:G:H8	1.81	0.45
23:BB:1340:U:H5'	50:BT:61:LEU:CD2	2.46	0.45
23:BB:1370:C:H2'	23:BB:1371:G:O4'	2.16	0.45
23:BB:1428:C:H2'	23:BB:1569:A:OP2	2.16	0.45
23:BB:1764:C:O2'	23:BB:1765:U:H5'	2.17	0.45
23:BB:1919:A:H3'	23:BB:1920:C:H6	1.81	0.45
23:BB:2000:C:O2'	23:BB:2001:C:H5'	2.17	0.45
23:BB:2142:A:H2'	23:BB:2143:C:O4'	2.16	0.45
23:BB:2351:G:HO2'	23:BB:2352:A:H8	1.63	0.45
23:BB:2415:G:H2'	23:BB:2416:C:C6	2.51	0.45
23:BB:2448:A:H4'	23:BB:2449:U:OP2	2.17	0.45
23:BB:2655:G:N2	23:BB:2664:G:H2'	2.31	0.45
23:BB:2708:G:O2'	23:BB:2709:G:H5'	2.17	0.45
23:BB:419:U:H2'	23:BB:420:C:H6	1.80	0.45
23:BB:680:C:H2'	23:BB:681:G:C8	2.51	0.45
23:BB:858:G:H21	23:BB:2268:A:C3'	2.26	0.45
23:BB:903:C:H2'	23:BB:904:G:C8	2.51	0.45
29:BE:17:THR:HG22	29:BE:106:LYS:HE2	1.99	0.45
48:BG:104:LEU:HB2	48:BG:112:VAL:HB	1.97	0.45
24:BI:29:GLN:HA	24:BI:29:GLN:NE2	2.31	0.45
41:BJ:7:LYS:C	41:BJ:9:GLU:H	2.19	0.45
38:BM:29:GLY:CA	38:BM:106:ASP:HB2	2.47	0.45
42:BN:28:LEU:HD23	42:BN:34:ILE:HG13	1.98	0.45
23:BB:2881:U:O3'	42:BN:96:ARG:HG3	2.16	0.45
43:BO:30:ARG:H	43:BO:97:PHE:HE2	1.62	0.45
49:BR:2:TYR:CD1	49:BR:42:ALA:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:38:VAL:HB	49:BR:59:ILE:HD11	1.99	0.45
51:BZ:32:ASN:O	51:BZ:33:LEU:C	2.54	0.45
1:CA:1122:U:H2'	1:CA:1123:U:H6	1.80	0.45
1:CA:1218:C:H2'	1:CA:1219:A:H8	1.79	0.45
1:CA:1492:A:H61	1:CA:1494:G:H1'	1.82	0.45
1:CA:35:G:O2'	1:CA:36:C:H5'	2.16	0.45
1:CA:510:A:N3	1:CA:543:U:H1'	2.31	0.45
1:CA:730:G:O2'	1:CA:766:A:H5'	2.15	0.45
2:CC:133:MET:CE	2:CC:165:GLU:HG3	2.47	0.45
2:CC:88:LYS:HD3	2:CC:88:LYS:C	2.36	0.45
3:CD:146:GLU:CD	3:CD:146:GLU:H	2.19	0.45
3:CD:144:ILE:HD13	3:CD:154:VAL:HG11	1.98	0.45
3:CD:54:LEU:CD1	3:CD:55:ARG:HH21	2.30	0.45
4:CE:114:LEU:HD13	4:CE:122:VAL:HG21	1.98	0.45
5:CF:17:GLN:HB3	5:CF:17:GLN:HE21	1.59	0.45
5:CF:64:VAL:HG12	5:CF:65:GLU:N	2.32	0.45
6:CG:26:VAL:HA	6:CG:42:VAL:HG21	1.98	0.45
8:CI:85:ALA:O	8:CI:88:GLU:HG2	2.16	0.45
31:D0:42:ILE:HG22	31:D0:43:THR:O	2.17	0.45
36:D2:9:VAL:CG1	36:D2:10:LEU:N	2.76	0.45
34:D3:30:HIS:HD2	34:D3:31:ILE:H	1.65	0.45
22:DA:9:G:P	43:DO:25:ARG:HH12	2.39	0.45
23:DB:1064:C:H2'	23:DB:1065:U:O4'	2.16	0.45
23:DB:2282:G:H4'	23:DB:2283:C:O5'	2.16	0.45
23:DB:2581:G:N3	23:DB:2581:G:H2'	2.30	0.45
23:DB:557:C:H2'	23:DB:558:U:H6	1.81	0.45
23:DB:740:C:O2'	23:DB:741:U:H5'	2.17	0.45
25:DC:239:PHE:C	25:DC:241:LYS:H	2.18	0.45
23:DB:2032:G:N2	26:DD:150:GLN:HB3	2.31	0.45
29:DE:15:SER:HB3	29:DE:18:THR:OG1	2.16	0.45
41:DJ:93:ILE:CA	41:DJ:97:PRO:HG3	2.45	0.45
38:DM:101:VAL:HG22	38:DM:101:VAL:O	2.17	0.45
42:DN:48:VAL:HG23	42:DN:49:GLU:N	2.32	0.45
42:DN:55:ALA:HA	42:DN:80:PHE:CE1	2.50	0.45
43:DO:105:ALA:C	43:DO:107:ALA:H	2.19	0.45
28:DP:47:ILE:HG13	28:DP:48:ALA:H	1.82	0.45
44:DQ:7:VAL:HG23	44:DQ:8:ILE:N	2.31	0.45
45:DS:14:ALA:C	45:DS:16:LYS:H	2.19	0.45
51:DZ:27:ARG:O	51:DZ:28:ARG:HB3	2.15	0.45
1:AA:1316:G:H22	1:AA:1318:A:H3'	1.79	0.45
1:AA:309:A:O2'	1:AA:310:G:H5'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:473:U:OP1	13:AP:76:LYS:HE2	2.15	0.45
1:AA:486:U:H2'	1:AA:487:A:H8	1.81	0.45
1:AA:51:A:H4'	1:AA:52:C:OP2	2.15	0.45
1:AA:761:G:H2'	1:AA:762:U:C6	2.52	0.45
18:AB:83:ALA:HA	18:AB:88:GLN:HE21	1.80	0.45
2:AC:63:ILE:HB	2:AC:98:ALA:CB	2.46	0.45
4:AE:152:VAL:O	4:AE:156:ARG:HG2	2.16	0.45
8:AI:123:ARG:HH11	8:AI:123:ARG:CB	2.25	0.45
11:AL:13:ARG:HB2	11:AL:14:LYS:H	1.50	0.45
21:AN:26:LEU:O	21:AN:29:ILE:HB	2.16	0.45
1:AA:981:U:O3'	21:AN:60:ARG:HD3	2.16	0.45
14:AQ:30:HIS:CG	14:AQ:33:TYR:HB2	2.51	0.45
16:AS:33:TRP:CD2	16:AS:33:TRP:N	2.83	0.45
23:BB:1183:U:H2'	23:BB:1184:U:C6	2.50	0.45
23:BB:1212:G:H1'	23:BB:1236:G:N2	2.32	0.45
23:BB:1238:G:O2'	23:BB:1239:G:H5'	2.17	0.45
23:BB:1316:U:H2'	23:BB:1317:G:H8	1.81	0.45
23:BB:1864:U:O2'	23:BB:1865:U:H5'	2.16	0.45
23:BB:2581:G:C8	23:BB:2610:C:N4	2.85	0.45
23:BB:2822:G:H2'	23:BB:2823:A:H5''	1.97	0.45
23:BB:341:C:H2'	23:BB:342:A:H8	1.81	0.45
23:BB:359:G:C2'	23:BB:360:U:H5'	2.47	0.45
23:BB:464:U:H2'	23:BB:465:G:O4'	2.15	0.45
23:BB:944:C:H5'	23:BB:945:A:C5'	2.46	0.45
23:BB:945:A:H3'	23:BB:946:C:H5''	1.98	0.45
25:BC:5:CYS:HB2	25:BC:15:VAL:O	2.17	0.45
25:BC:201:LEU:O	25:BC:202:ARG:O	2.34	0.45
48:BG:15:ASP:O	48:BG:16:VAL:HB	2.16	0.45
22:BA:49:C:OP1	43:BO:101:GLY:HA3	2.17	0.45
35:BV:63:ILE:HB	35:BV:70:ILE:HG12	1.98	0.45
35:BV:35:GLU:HB3	35:BV:93:ARG:CZ	2.47	0.45
23:BB:2330:G:H1'	52:BW:38:ARG:HB2	1.97	0.45
30:BY:47:ILE:HG21	30:BY:56:VAL:CG2	2.47	0.45
1:CA:1001:C:H2'	1:CA:1002:G:C8	2.51	0.45
1:CA:976:G:N2	1:CA:1363:A:C2	2.81	0.45
1:CA:286:C:H2'	1:CA:287:U:H6	1.81	0.45
1:CA:386:C:O2'	1:CA:387:U:H5'	2.16	0.45
1:CA:647:C:H2'	1:CA:648:A:H8	1.81	0.45
1:CA:884:U:H4'	1:CA:885:G:H5''	1.99	0.45
1:CA:989:U:C2'	1:CA:990:C:H5'	2.46	0.45
18:CB:161:PHE:HD2	18:CB:183:PHE:HB2	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:85:ILE:HG22	5:CF:86:ARG:N	2.31	0.45
8:CI:11:ARG:HH22	8:CI:108:ARG:HB2	1.80	0.45
8:CI:35:GLU:HA	8:CI:39:GLY:HA2	1.97	0.45
12:CM:36:ALA:HB2	12:CM:58:GLU:OE1	2.16	0.45
1:CA:1320:C:H1'	16:CS:72:GLU:HB3	1.99	0.45
23:DB:101:A:C2'	23:DB:102:U:H5''	2.46	0.45
23:DB:1055:G:H2'	23:DB:1056:G:O4'	2.17	0.45
23:DB:1183:U:H2'	23:DB:1184:U:C6	2.51	0.45
23:DB:1745:A:H2'	23:DB:1746:A:O4'	2.17	0.45
23:DB:1820:U:H3	25:DC:197:ALA:HA	1.82	0.45
23:DB:1902:C:H2'	23:DB:1903:G:O4'	2.17	0.45
23:DB:2379:G:C5'	43:DO:21:LEU:HD11	2.47	0.45
23:DB:2699:C:H2'	23:DB:2700:A:H8	1.81	0.45
23:DB:2722:G:H2'	23:DB:2723:C:H6	1.81	0.45
23:DB:2742:G:H2'	23:DB:2743:U:H6	1.80	0.45
29:DE:161:ALA:HA	29:DE:164:LEU:HB2	1.99	0.45
29:DE:112:LEU:HD11	29:DE:180:LEU:HB3	1.98	0.45
47:DF:110:ILE:HG22	47:DF:113:PHE:HB3	1.99	0.45
47:DF:168:LEU:C	47:DF:170:ALA:H	2.19	0.45
47:DF:11:VAL:HG21	47:DF:172:PHE:CE1	2.52	0.45
48:DG:123:GLU:O	48:DG:125:PRO:HD3	2.16	0.45
40:DH:124:THR:O	40:DH:128:HIS:HE1	1.98	0.45
24:DI:91:LYS:HD2	24:DI:91:LYS:N	2.31	0.45
27:DK:41:ILE:HG13	27:DK:42:THR:N	2.31	0.45
27:DK:68:GLY:CA	27:DK:78:ARG:HB3	2.47	0.45
37:DL:69:ARG:HE	37:DL:69:ARG:HA	1.80	0.45
38:DM:114:ARG:HG3	38:DM:130:PHE:CD1	2.51	0.45
38:DM:71:LYS:O	38:DM:92:TRP:HA	2.15	0.45
28:DP:20:ARG:HB3	28:DP:23:ASP:OD2	2.17	0.45
27:DK:79:PHE:HD2	28:DP:69:VAL:HG12	1.82	0.45
44:DQ:75:TYR:O	44:DQ:79:ILE:HG22	2.17	0.45
49:DR:75:VAL:O	49:DR:76:LYS:HG3	2.16	0.45
1:AA:1026:G:H8	1:AA:1026:G:O5'	2.00	0.45
1:AA:1060:U:H2'	1:AA:1061:G:H8	1.80	0.45
1:AA:1113:C:H2'	1:AA:1114:C:H6	1.82	0.45
1:AA:1207:G:H2'	1:AA:1208:C:H6	1.82	0.45
1:AA:1306:A:N6	1:AA:1331:G:H1'	2.32	0.45
1:AA:1409:C:O2	1:AA:1409:C:H2'	2.16	0.45
1:AA:1448:C:O2'	1:AA:1449:C:H5'	2.16	0.45
1:AA:1409:C:N3	1:AA:1492:A:C2	2.83	0.45
1:AA:47:C:H4'	1:AA:48:C:O5'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:126:G:C4'	1:AA:634:C:H1'	2.47	0.45
18:AB:65:LYS:HZ2	18:AB:89:PHE:HE1	1.64	0.45
2:AC:10:ARG:HH11	2:AC:10:ARG:HG3	1.81	0.45
2:AC:26:LYS:HG3	2:AC:27:GLU:HG3	1.98	0.45
3:AD:24:VAL:HG12	3:AD:160:LEU:HB3	1.98	0.45
3:AD:138:PRO:HA	3:AD:181:PHE:CD2	2.52	0.45
6:AG:50:ALA:HB2	6:AG:57:GLU:CB	2.46	0.45
9:AJ:67:ILE:HG23	21:AN:94:GLY:O	2.17	0.45
12:AM:11:HIS:HA	12:AM:43:LYS:HD3	1.98	0.45
21:AN:44:VAL:O	21:AN:48:GLN:HB3	2.17	0.45
1:AA:564:C:N1	14:AQ:32:ILE:HD11	2.31	0.45
14:AQ:55:GLY:O	14:AQ:81:ALA:HB2	2.16	0.45
15:AR:28:LEU:C	15:AR:30:ASN:H	2.18	0.45
31:B0:27:LEU:HD21	45:BS:38:TYR:CE1	2.51	0.45
36:B2:22:MET:HG2	36:B2:31:LEU:HD12	1.99	0.45
23:BB:1179:G:H2'	23:BB:1180:U:O4'	2.16	0.45
23:BB:1383:A:H2	23:BB:1405:U:O2	1.99	0.45
23:BB:1454:C:C5	42:BN:64:ARG:HG2	2.52	0.45
23:BB:1537:G:H5''	23:BB:1537:G:N3	2.31	0.45
23:BB:1788:C:H2'	23:BB:1789:A:H8	1.80	0.45
23:BB:182:A:H2'	23:BB:183:C:H6	1.81	0.45
23:BB:2037:A:H2'	23:BB:2038:G:H8	1.82	0.45
23:BB:2484:G:O2'	23:BB:2485:G:H5'	2.17	0.45
23:BB:1664:A:H1'	23:BB:2726:A:N1	2.32	0.45
23:BB:728:G:O2'	23:BB:730:A:H8	1.99	0.45
23:BB:863:A:O2'	23:BB:864:G:H5'	2.17	0.45
23:BB:878:A:N3	23:BB:878:A:C2'	2.78	0.45
25:BC:14:HIS:O	25:BC:16:VAL:HG23	2.17	0.45
25:BC:211:ARG:C	25:BC:213:ARG:H	2.19	0.45
25:BC:226:PRO:HA	25:BC:232:GLY:HA3	1.99	0.45
29:BE:58:LYS:N	29:BE:58:LYS:HD3	2.32	0.45
29:BE:98:LYS:HE3	29:BE:99:LYS:HG2	1.98	0.45
47:BF:135:ILE:C	47:BF:137:PHE:H	2.19	0.45
40:BH:58:LEU:C	40:BH:60:GLU:H	2.20	0.45
27:BK:20:MET:C	27:BK:41:ILE:HG13	2.36	0.45
23:BB:662:G:O3'	37:BL:16:GLY:HA2	2.17	0.45
37:BL:68:SER:CB	37:BL:71:ALA:HB3	2.45	0.45
42:BN:101:GLY:CA	42:BN:109:PRO:HA	2.44	0.45
28:BP:89:GLY:H	28:BP:112:ARG:NH1	2.14	0.45
28:BP:21:PRO:CG	28:BP:91:VAL:HG21	2.37	0.45
44:BQ:57:ARG:C	44:BQ:59:LEU:H	2.18	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:75:VAL:O	49:BR:76:LYS:HG3	2.17	0.45
50:BT:23:ALA:C	50:BT:25:GLU:H	2.20	0.45
46:BU:35:VAL:HB	46:BU:38:ILE:HG21	1.99	0.45
35:BV:14:LYS:O	35:BV:18:ARG:HB2	2.17	0.45
35:BV:65:VAL:O	35:BV:67:GLY:N	2.47	0.45
52:BW:30:VAL:HA	52:BW:60:ALA:O	2.16	0.45
1:CA:1010:U:O2'	1:CA:1011:C:H5'	2.17	0.45
1:CA:1128:C:H1'	1:CA:1148:U:O4	2.15	0.45
1:CA:1249:C:H2'	1:CA:1250:A:H5'	1.97	0.45
1:CA:976:G:N1	1:CA:1362:A:H3'	2.32	0.45
1:CA:168:G:O2'	1:CA:169:C:H5'	2.16	0.45
1:CA:178:C:H2'	1:CA:179:A:H8	1.81	0.45
1:CA:45:G:H2'	1:CA:46:G:C8	2.52	0.45
1:CA:989:U:O2'	1:CA:990:C:H5'	2.16	0.45
18:CB:101:THR:HG22	18:CB:174:GLU:CD	2.36	0.45
18:CB:151:LYS:HG3	18:CB:152:ASP:H	1.80	0.45
18:CB:169:HIS:HA	18:CB:172:ILE:CD1	2.46	0.45
2:CC:106:ARG:HD3	2:CC:107:LYS:NZ	2.31	0.45
6:CG:19:SER:HB3	6:CG:22:LEU:CB	2.40	0.45
8:CI:30:ASN:O	8:CI:31:GLN:HB2	2.17	0.45
12:CM:65:GLU:N	12:CM:65:GLU:OE2	2.49	0.45
20:CO:10:LYS:HG3	20:CO:11:ILE:N	2.32	0.45
13:CP:67:ILE:CG1	13:CP:72:ALA:HB2	2.44	0.45
34:D3:40:LYS:O	34:D3:43:LEU:N	2.47	0.45
22:DA:94:A:H2'	22:DA:95:U:O4'	2.16	0.45
23:DB:1169:A:H2'	23:DB:1170:C:C6	2.51	0.45
23:DB:1172:C:H2'	23:DB:1173:U:O4'	2.17	0.45
23:DB:1228:G:H2'	23:DB:1229:C:C6	2.52	0.45
23:DB:1418:G:H1'	23:DB:1580:A:H61	1.82	0.45
23:DB:1535:A:O2'	23:DB:1536:C:H5'	2.16	0.45
23:DB:1538:G:O2'	23:DB:1539:U:H5'	2.17	0.45
23:DB:1654:A:H2'	23:DB:1655:A:C8	2.51	0.45
23:DB:1745:A:H2'	23:DB:1746:A:H8	1.80	0.45
23:DB:1827:U:O2'	23:DB:1828:G:H5'	2.16	0.45
23:DB:2527:C:H2'	23:DB:2528:U:H6	1.81	0.45
23:DB:2533:U:H2'	23:DB:2534:A:O4'	2.15	0.45
23:DB:2732:G:H5'	23:DB:2733:A:O4'	2.16	0.45
23:DB:483:A:H2'	23:DB:484:C:O4'	2.17	0.45
23:DB:552:U:O2'	23:DB:553:G:H5'	2.17	0.45
23:DB:55:G:H2'	23:DB:56:A:H8	1.82	0.45
25:DC:183:VAL:HG22	25:DC:184:GLU:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:184:GLU:O	25:DC:186:ASP:N	2.42	0.45
25:DC:270:ARG:HH11	25:DC:270:ARG:HB3	1.82	0.45
29:DE:16:GLU:O	29:DE:20:GLY:HA3	2.17	0.45
23:DB:588:U:H1'	29:DE:85:PHE:CD2	2.51	0.45
47:DF:31:GLU:HB3	47:DF:157:THR:HA	1.98	0.45
47:DF:121:PHE:CE1	47:DF:162:ASP:HB2	2.51	0.45
48:DG:1:SER:H1	48:DG:61:TRP:HE3	1.64	0.45
40:DH:133:GLN:HG2	40:DH:139:PHE:HB3	1.98	0.45
40:DH:30:LEU:O	40:DH:35:LYS:HB2	2.15	0.45
27:DK:19:VAL:CG1	27:DK:41:ILE:HG12	2.42	0.45
27:DK:18:ARG:HB2	27:DK:45:GLU:HG2	1.97	0.45
27:DK:71:ARG:NE	27:DK:72:PRO:HD3	2.30	0.45
27:DK:79:PHE:CD1	27:DK:79:PHE:N	2.84	0.45
42:DN:24:MET:CG	42:DN:44:LEU:HD22	2.46	0.45
42:DN:96:ARG:N	42:DN:96:ARG:HD3	2.31	0.45
27:DK:108:ARG:NH1	28:DP:34:GLY:HA2	2.32	0.45
28:DP:62:LYS:HB3	28:DP:69:VAL:HG22	1.99	0.45
44:DQ:30:VAL:HG11	44:DQ:33:VAL:HG22	1.98	0.45
41:DJ:44:TYR:HB2	44:DQ:63:ARG:HG2	1.97	0.45
35:DV:11:GLU:CD	35:DV:16:ALA:HB1	2.37	0.45
1:AA:1130:A:N6	1:AA:1143:G:N2	2.64	0.45
1:AA:1307:U:H2'	1:AA:1308:U:H6	1.81	0.45
1:AA:197:A:H4'	1:AA:198:G:O5'	2.17	0.45
1:AA:410:G:H1'	1:AA:432:A:H61	1.82	0.45
1:AA:410:G:H8	1:AA:410:G:O5'	1.99	0.45
1:AA:424:G:O2'	1:AA:425:G:H5'	2.16	0.45
1:AA:640:A:O2'	1:AA:641:U:H5'	2.17	0.45
1:AA:981:U:H2'	1:AA:982:U:C5	2.52	0.45
1:AA:986:U:H2'	1:AA:987:G:O4'	2.17	0.45
18:AB:53:LEU:C	18:AB:55:GLU:H	2.19	0.45
2:AC:131:ARG:HB3	2:AC:135:ARG:NH2	2.32	0.45
7:AH:118:ALA:HB3	7:AH:120:LEU:HD22	1.99	0.45
8:AI:112:ARG:NH2	8:AI:114:LYS:HA	2.32	0.45
8:AI:123:ARG:HB2	8:AI:123:ARG:NH1	2.24	0.45
8:AI:33:SER:HB3	8:AI:36:GLN:HB2	1.99	0.45
10:AK:125:LYS:O	19:AU:33:ARG:NH1	2.50	0.45
11:AL:28:GLN:HB2	11:AL:80:LEU:HG	1.99	0.45
11:AL:36:VAL:O	11:AL:36:VAL:HG23	2.16	0.45
13:AP:72:ALA:HA	13:AP:75:ILE:HD12	1.97	0.45
15:AR:52:ARG:O	15:AR:56:ARG:HG3	2.17	0.45
17:AT:78:LEU:O	17:AT:82:ILE:HG23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:B0:39:ARG:O	31:B0:40:HIS:HB2	2.17	0.45
33:B1:32:LYS:NZ	33:B1:52:LYS:HG2	2.32	0.45
34:B3:56:LEU:O	34:B3:59:ALA:HB3	2.15	0.45
22:BA:9:G:O2'	22:BA:10:G:H5'	2.17	0.45
23:BB:1105:U:H2'	23:BB:1106:G:C8	2.52	0.45
23:BB:1241:A:O4'	23:BB:1241:A:N3	2.49	0.45
23:BB:2032:G:N2	26:BD:150:GLN:HB3	2.32	0.45
23:BB:2897:U:H2'	23:BB:2898:U:C6	2.51	0.45
23:BB:68:G:H2'	23:BB:69:C:C6	2.52	0.45
23:BB:919:U:H6	23:BB:919:U:O5'	1.99	0.45
47:BF:11:VAL:HG21	47:BF:172:PHE:CE1	2.52	0.45
47:BF:127:TYR:CE2	47:BF:176:PHE:HE2	2.35	0.45
47:BF:1:ALA:N	47:BF:2:LYS:HE2	2.31	0.45
48:BG:17:LYS:O	48:BG:23:ILE:HG23	2.16	0.45
40:BH:6:LEU:HD12	40:BH:6:LEU:N	2.32	0.45
41:BJ:102:GLU:HA	41:BJ:105:VAL:HG22	1.98	0.45
27:BK:72:PRO:O	27:BK:74:GLY:N	2.50	0.45
27:BK:93:GLN:HB3	27:BK:94:PRO:HD2	1.98	0.45
37:BL:132:ARG:O	37:BL:136:GLU:HB2	2.16	0.45
38:BM:33:LEU:HD22	38:BM:128:THR:CB	2.43	0.45
38:BM:38:ARG:HG2	38:BM:98:PRO:HD3	1.99	0.45
43:BO:106:LEU:HG	43:BO:107:ALA:N	2.30	0.45
28:BP:109:ILE:HG13	28:BP:109:ILE:O	2.17	0.45
49:BR:31:GLU:H	49:BR:63:VAL:CG2	2.30	0.45
45:BS:49:LYS:C	45:BS:51:LEU:H	2.19	0.45
50:BT:30:ILE:O	50:BT:85:VAL:HG23	2.16	0.45
50:BT:32:LEU:N	50:BT:83:ALA:CB	2.79	0.45
50:BT:53:VAL:CG1	50:BT:87:LEU:HB3	2.46	0.45
1:CA:1001:C:H2'	1:CA:1002:G:H8	1.81	0.45
1:CA:1015:G:H2'	1:CA:1016:A:H8	1.82	0.45
1:CA:144:G:H2'	1:CA:145:G:O4'	2.17	0.45
1:CA:1486:G:H2'	1:CA:1487:G:C8	2.51	0.45
1:CA:337:G:O2'	1:CA:338:A:H5'	2.17	0.45
1:CA:764:C:N4	1:CA:812:G:H1	2.15	0.45
18:CB:103:TRP:CZ3	18:CB:107:ARG:HB3	2.52	0.45
18:CB:116:LEU:HD23	18:CB:119:GLN:CG	2.46	0.45
2:CC:100:ILE:HG23	2:CC:100:ILE:O	2.17	0.45
2:CC:172:VAL:HG12	2:CC:174:LEU:CD1	2.47	0.45
2:CC:38:VAL:O	2:CC:41:TYR:HB3	2.17	0.45
3:CD:155:LYS:HA	3:CD:158:LEU:HD12	1.98	0.45
31:D0:16:ARG:HA	31:D0:19:ASP:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:D0:47:TYR:CD1	31:D0:47:TYR:N	2.85	0.45
23:DB:1117:C:H2'	23:DB:1118:C:C6	2.52	0.45
23:DB:1169:A:H2'	23:DB:1170:C:H6	1.82	0.45
23:DB:1322:A:C2'	23:DB:1323:C:H5'	2.47	0.45
23:DB:1461:C:O2'	23:DB:1462:C:H5'	2.16	0.45
23:DB:1684:G:H2'	23:DB:1685:C:H6	1.81	0.45
23:DB:1719:G:O2'	23:DB:1720:U:H5'	2.17	0.45
23:DB:172:A:H2'	23:DB:173:A:C8	2.51	0.45
23:DB:1867:G:O2'	23:DB:1868:C:H5'	2.16	0.45
23:DB:1951:U:O2	23:DB:1953:A:H3'	2.17	0.45
23:DB:2201:G:H2'	23:DB:2202:U:C6	2.51	0.45
23:DB:2361:G:O2'	23:DB:2362:C:H5'	2.17	0.45
23:DB:2448:A:H4'	23:DB:2449:U:OP2	2.16	0.45
23:DB:2599:G:O2'	23:DB:2600:A:H5'	2.16	0.45
23:DB:2721:A:H2'	23:DB:2722:G:C8	2.51	0.45
23:DB:360:U:H2'	23:DB:361:G:N9	2.31	0.45
23:DB:988:A:OP1	30:DY:11:SER:HB3	2.16	0.45
25:DC:226:PRO:HA	25:DC:232:GLY:HA3	1.98	0.45
25:DC:63:ILE:HD13	25:DC:63:ILE:HA	1.84	0.45
25:DC:93:VAL:HG13	25:DC:101:ARG:H	1.80	0.45
26:DD:68:PHE:CB	26:DD:73:VAL:HG23	2.47	0.45
26:DD:8:LYS:HB2	26:DD:201:LEU:HD21	1.99	0.45
26:DD:92:VAL:O	26:DD:94:GLN:N	2.50	0.45
29:DE:17:THR:HG22	29:DE:106:LYS:HE2	1.98	0.45
47:DF:41:GLU:O	47:DF:43:ILE:N	2.49	0.45
47:DF:42:ALA:O	47:DF:46:LYS:HG3	2.16	0.45
48:DG:77:GLY:HA3	48:DG:135:ALA:O	2.16	0.45
24:DI:45:THR:HA	24:DI:48:ILE:CG2	2.40	0.45
41:DJ:3:THR:HB	41:DJ:44:TYR:CE1	2.52	0.45
27:DK:51:LYS:O	27:DK:52:VAL:HG13	2.16	0.45
37:DL:10:GLU:C	37:DL:12:SER:H	2.20	0.45
44:DQ:71:ASN:HD22	44:DQ:109:VAL:HG21	1.81	0.45
44:DQ:9:ALA:O	44:DQ:11:ALA:N	2.49	0.45
49:DR:74:ILE:HB	49:DR:87:GLN:O	2.16	0.45
45:DS:3:THR:HB	45:DS:62:ASP:CB	2.47	0.45
46:DU:86:PHE:CE1	46:DU:88:ASP:HB2	2.51	0.45
52:DW:24:ARG:CD	52:DW:65:LYS:HG2	2.47	0.45
52:DW:76:ARG:HH21	52:DW:76:ARG:HA	1.79	0.45
51:DZ:77:LYS:O	51:DZ:78:TYR:HB3	2.15	0.45
1:AA:1117:A:H4'	8:AI:105:ARG:HH12	1.82	0.45
1:AA:114:U:O2'	1:AA:115:G:H5'	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1203:C:H2'	1:AA:1204:A:H8	1.81	0.45
1:AA:1240:U:C3'	1:AA:1241:G:H5'	2.46	0.45
1:AA:728:A:H2'	1:AA:729:A:C8	2.51	0.45
18:AB:119:GLN:C	18:AB:121:GLN:H	2.19	0.45
2:AC:139:ASN:HD22	2:AC:139:ASN:HA	1.55	0.45
2:AC:185:THR:HB	2:AC:197:VAL:O	2.17	0.45
2:AC:205:GLU:O	2:AC:206:ILE:C	2.53	0.45
3:AD:35:GLN:O	3:AD:37:PRO:HD3	2.16	0.45
5:AF:38:ARG:O	5:AF:62:MET:O	2.35	0.45
6:AG:103:ILE:HG21	6:AG:123:LEU:HD23	1.98	0.45
6:AG:68:VAL:HA	6:AG:137:ARG:HG3	1.97	0.45
6:AG:14:ASP:OD1	6:AG:17:PHE:HB2	2.17	0.45
11:AL:17:LYS:H	11:AL:17:LYS:HE3	1.82	0.45
12:AM:86:ARG:HB3	12:AM:86:ARG:HE	1.56	0.45
21:AN:14:ALA:HB1	21:AN:18:LYS:HE3	1.98	0.45
14:AQ:8:GLN:HA	14:AQ:58:VAL:O	2.17	0.45
16:AS:14:LEU:HD13	16:AS:34:SER:OG	2.17	0.45
19:AU:13:VAL:O	19:AU:13:VAL:HG13	2.17	0.45
23:BB:1064:C:H2'	23:BB:1065:U:O4'	2.17	0.45
23:BB:2266:A:H4'	23:BB:2267:A:O5'	2.17	0.45
23:BB:2331:G:O2'	23:BB:2332:C:H5'	2.16	0.45
23:BB:2521:C:H2'	23:BB:2522:U:C6	2.52	0.45
23:BB:2527:C:O2'	23:BB:2528:U:H5'	2.16	0.45
23:BB:264:C:H2'	23:BB:265:A:C5'	2.45	0.45
23:BB:2867:G:N3	23:BB:2867:G:C2'	2.78	0.45
23:BB:302:C:H2'	23:BB:303:G:C8	2.50	0.45
23:BB:738:G:H2'	23:BB:739:A:C8	2.52	0.45
23:BB:875:G:H2'	23:BB:876:C:O2	2.17	0.45
26:BD:35:THR:O	26:BD:71:ALA:HB1	2.17	0.45
26:BD:55:LYS:C	26:BD:57:ALA:H	2.19	0.45
29:BE:40:ARG:NH2	29:BE:92:HIS:NE2	2.65	0.45
47:BF:106:ALA:HA	47:BF:135:ILE:CD1	2.46	0.45
47:BF:134:GLN:HE21	47:BF:134:GLN:HB3	1.51	0.45
47:BF:147:ARG:HD2	47:BF:147:ARG:O	2.17	0.45
47:BF:42:ALA:O	47:BF:46:LYS:HG3	2.17	0.45
47:BF:73:VAL:O	47:BF:74:ALA:HB2	2.16	0.45
24:BI:116:MET:SD	24:BI:124:MET:HB2	2.56	0.45
23:BB:1952:A:OP1	27:BK:42:THR:HG21	2.16	0.45
23:BB:632:A:H1'	37:BL:66:PHE:HE2	1.82	0.45
42:BN:101:GLY:O	42:BN:102:PHE:HB2	2.16	0.45
28:BP:58:PHE:HD1	28:BP:73:PHE:HD2	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BJ:44:TYR:HB2	44:BQ:63:ARG:HG2	1.97	0.45
44:BQ:60:TRP:HB3	44:BQ:92:LYS:O	2.17	0.45
45:BS:49:LYS:C	45:BS:51:LEU:N	2.69	0.45
23:BB:327:G:N2	46:BU:67:SER:HB2	2.31	0.45
30:BY:4:ILE:HG12	30:BY:38:GLU:O	2.17	0.45
1:CA:1157:A:H4'	1:CA:1158:C:O5'	2.17	0.45
1:CA:537:G:H2'	1:CA:538:G:C8	2.52	0.45
1:CA:761:G:H2'	1:CA:762:U:C6	2.52	0.45
2:CC:116:ALA:O	2:CC:117:ASP:C	2.54	0.45
3:CD:116:LEU:HD21	3:CD:153:ARG:HD3	1.97	0.45
3:CD:154:VAL:O	3:CD:157:ALA:HB3	2.16	0.45
4:CE:40:ASP:OD2	4:CE:42:ASN:HB3	2.17	0.45
7:CH:14:ARG:HE	7:CH:75:GLN:NE2	2.14	0.45
7:CH:82:LEU:O	7:CH:82:LEU:HD13	2.17	0.45
12:CM:90:HIS:HA	12:CM:108:ARG:HH21	1.79	0.45
20:CO:36:ILE:CD1	20:CO:59:MET:HG3	2.47	0.45
20:CO:8:THR:O	20:CO:11:ILE:HG13	2.16	0.45
13:CP:19:VAL:H	13:CP:38:PHE:HA	1.81	0.45
14:CQ:8:GLN:HB3	14:CQ:59:GLU:HG3	1.98	0.45
17:CT:4:LYS:HE3	17:CT:6:ALA:HB2	1.98	0.45
31:D0:28:SER:HB3	31:D0:37:HIS:HE1	1.81	0.45
33:D1:9:LYS:O	33:D1:50:GLU:HG3	2.17	0.45
36:D2:4:THR:O	36:D2:5:PHE:HB2	2.15	0.45
34:D3:7:ARG:NH1	34:D3:7:ARG:HG3	2.32	0.45
22:DA:17:C:O2'	22:DA:18:G:H5'	2.16	0.45
23:DB:1312:U:O2	23:DB:1603:A:N1	2.50	0.45
23:DB:1435:G:H2'	23:DB:1436:G:H8	1.82	0.45
23:DB:1812:U:H4'	25:DC:44:ASN:OD1	2.16	0.45
23:DB:2223:G:C2'	23:DB:2224:G:H5'	2.47	0.45
23:DB:2236:U:O2'	23:DB:2237:G:H5'	2.17	0.45
23:DB:2336:A:H61	52:DW:40:ARG:CG	2.28	0.45
23:DB:233:A:N6	23:DB:428:A:H61	2.14	0.45
23:DB:241:A:H8	23:DB:241:A:OP1	2.00	0.45
23:DB:2547:A:H2'	23:DB:2548:U:H6	1.79	0.45
23:DB:2720:U:H5''	28:DP:52:ARG:HH22	1.79	0.45
23:DB:409:G:H2'	23:DB:410:G:H8	1.80	0.45
23:DB:625:G:O2'	23:DB:626:A:H5'	2.17	0.45
23:DB:802:A:H2'	23:DB:803:U:H6	1.79	0.45
23:DB:902:C:H2'	23:DB:903:C:C6	2.52	0.45
23:DB:915:C:H3'	23:DB:916:G:C8	2.50	0.45
29:DE:114:ARG:HG3	29:DE:114:ARG:NH1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:110:ILE:HA	47:DF:111:ARG:NH2	2.31	0.45
48:DG:40:VAL:HG13	48:DG:64:ALA:CB	2.46	0.45
48:DG:16:VAL:HG11	48:DG:44:HIS:NE2	2.32	0.45
40:DH:119:ASN:O	40:DH:121:VAL:HG22	2.16	0.45
40:DH:31:VAL:CB	40:DH:32:PRO:CD	2.88	0.45
24:DI:1:ALA:C	24:DI:2:LYS:HD2	2.36	0.45
41:DJ:37:ARG:HE	41:DJ:110:PRO:HG3	1.82	0.45
37:DL:116:VAL:HG13	37:DL:117:THR:N	2.19	0.45
37:DL:124:GLY:N	37:DL:143:GLU:CG	2.72	0.45
38:DM:105:MET:HB2	38:DM:117:PHE:CZ	2.47	0.45
28:DP:20:ARG:CG	28:DP:21:PRO:HD2	2.42	0.45
44:DQ:73:ILE:HG21	44:DQ:109:VAL:HG13	1.98	0.45
49:DR:64:VAL:O	49:DR:65:ALA:HB3	2.17	0.45
46:DU:13:LEU:HD12	46:DU:13:LEU:N	2.31	0.45
46:DU:11:ILE:HG21	46:DU:78:LYS:CE	2.46	0.45
35:DV:29:ILE:HG13	35:DV:88:HIS:CE1	2.52	0.45
23:DB:929:U:H1'	30:DY:25:GLY:O	2.17	0.45
1:AA:1049:U:H5''	1:AA:1201:A:OP2	2.17	0.45
1:AA:1191:A:H2'	1:AA:1192:C:C6	2.51	0.45
1:AA:1257:A:N3	1:AA:1257:A:O4'	2.49	0.45
1:AA:1270:G:H8	1:AA:1270:G:O5'	2.00	0.45
1:AA:1325:C:O2'	1:AA:1326:U:H5'	2.17	0.45
1:AA:137:U:H2'	1:AA:138:G:C8	2.51	0.45
1:AA:1473:G:H2'	1:AA:1474:U:O4'	2.17	0.45
1:AA:473:U:H2'	1:AA:474:G:H8	1.81	0.45
18:AB:101:THR:HG23	18:AB:102:ASN:N	2.32	0.45
18:AB:163:ILE:CG2	18:AB:164:ASP:H	2.25	0.45
18:AB:83:ALA:CB	18:AB:90:PHE:HB3	2.46	0.45
4:AE:114:LEU:HD13	4:AE:122:VAL:HG21	1.97	0.45
4:AE:45:VAL:O	4:AE:71:ILE:HG22	2.17	0.45
5:AF:53:LYS:NZ	5:AF:53:LYS:N	2.64	0.45
6:AG:12:LEU:HA	6:AG:12:LEU:HD22	1.88	0.45
6:AG:46:LEU:HG	6:AG:57:GLU:HB2	1.99	0.45
11:AL:82:ARG:NH2	11:AL:95:HIS:ND1	2.65	0.45
20:AO:34:ALA:HB3	20:AO:35:GLN:OE1	2.16	0.45
13:AP:72:ALA:HA	13:AP:75:ILE:CD1	2.46	0.45
15:AR:37:LYS:HG2	15:AR:37:LYS:H	1.61	0.45
15:AR:70:THR:HB	15:AR:72:ARG:NH1	2.32	0.45
16:AS:18:VAL:HG13	16:AS:19:GLU:N	2.32	0.45
22:BA:96:G:O2'	22:BA:97:C:H5'	2.15	0.45
23:BB:1028:A:N3	23:BB:2486:C:O2'	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1196:C:H2'	23:BB:1197:G:H8	1.79	0.45
23:BB:1219:U:H2'	23:BB:1220:G:H8	1.82	0.45
23:BB:1374:G:H2'	23:BB:1375:U:H6	1.82	0.45
23:BB:1439:A:N3	23:BB:1553:A:C6	2.85	0.45
23:BB:1469:A:H2'	23:BB:1470:A:H8	1.81	0.45
23:BB:121:G:H1'	23:BB:148:U:N3	2.31	0.45
23:BB:1729:U:H5	23:BB:1731:G:H22	1.62	0.45
23:BB:1745:A:H2'	23:BB:1746:A:O4'	2.16	0.45
23:BB:2291:U:H2'	23:BB:2292:U:H6	1.81	0.45
23:BB:2511:U:H2'	23:BB:2512:C:C6	2.52	0.45
23:BB:2709:G:H2'	23:BB:2710:C:H6	1.82	0.45
23:BB:2732:G:H5'	23:BB:2733:A:O4'	2.16	0.45
23:BB:2888:C:H2'	23:BB:2889:C:H6	1.82	0.45
23:BB:383:C:N3	23:BB:391:A:N6	2.65	0.45
23:BB:596:U:O2'	23:BB:597:G:H5'	2.17	0.45
23:BB:657:U:H2'	23:BB:658:U:C6	2.52	0.45
23:BB:672:C:H2'	23:BB:673:C:C6	2.51	0.45
23:BB:735:A:N7	23:BB:761:A:H2	2.15	0.45
23:BB:811:U:O2	23:BB:1250:G:H2'	2.17	0.45
23:BB:838:C:O2'	23:BB:839:U:H5'	2.17	0.45
26:BD:117:GLY:HA2	26:BD:164:GLN:HE22	1.82	0.45
26:BD:195:GLY:HA3	26:BD:199:SER:HB2	1.98	0.45
26:BD:55:LYS:HZ3	26:BD:56:LYS:HG2	1.81	0.45
26:BD:58:ASN:OD1	26:BD:59:ARG:HG3	2.16	0.45
26:BD:7:LYS:HG3	26:BD:198:GLY:O	2.16	0.45
29:BE:29:HIS:CD2	37:BL:8:PRO:HA	2.52	0.45
47:BF:134:GLN:HB3	47:BF:149:ARG:HB3	1.99	0.45
47:BF:26:GLN:O	47:BF:27:VAL:C	2.54	0.45
47:BF:41:GLU:O	47:BF:43:ILE:N	2.50	0.45
48:BG:106:LEU:O	48:BG:108:PHE:HD1	1.99	0.45
48:BG:155:PRO:CA	48:BG:170:THR:HA	2.47	0.45
24:BI:107:GLU:HA	24:BI:110:GLN:OE1	2.17	0.45
41:BJ:42:ALA:O	41:BJ:44:TYR:N	2.50	0.45
37:BL:93:ASN:O	37:BL:95:LEU:HD12	2.17	0.45
38:BM:29:GLY:HA2	38:BM:106:ASP:HB2	1.97	0.45
42:BN:24:MET:CG	42:BN:44:LEU:HD22	2.47	0.45
42:BN:61:ALA:C	42:BN:63:ARG:N	2.69	0.45
31:B0:41:HIS:HB2	42:BN:99:LYS:O	2.17	0.45
28:BP:62:LYS:HB3	28:BP:69:VAL:CG2	2.46	0.45
44:BQ:9:ALA:C	44:BQ:11:ALA:N	2.70	0.45
49:BR:1:MET:C	49:BR:2:TYR:HD1	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:41:ILE:HD13	49:BR:41:ILE:N	2.31	0.45
49:BR:58:VAL:HG13	49:BR:59:ILE:N	2.31	0.45
45:BS:14:ALA:C	45:BS:16:LYS:H	2.19	0.45
50:BT:74:ILE:HG13	50:BT:75:GLY:N	2.30	0.45
52:BW:65:LYS:HZ2	52:BW:84:GLU:HB3	1.80	0.45
39:BX:17:GLU:CD	39:BX:18:LEU:N	2.70	0.45
39:BX:25:GLN:HE21	39:BX:25:GLN:HB3	1.54	0.45
30:BY:16:LEU:O	30:BY:19:HIS:HB2	2.15	0.45
51:BZ:27:ARG:O	51:BZ:28:ARG:HB3	2.17	0.45
1:CA:1251:A:N3	1:CA:1369:C:O2'	2.41	0.45
1:CA:1287:A:H2'	1:CA:1288:A:O4'	2.17	0.45
1:CA:32:A:H2'	1:CA:33:A:H8	1.81	0.45
1:CA:55:A:OP2	1:CA:352:C:N4	2.49	0.45
1:CA:203:G:H1'	1:CA:465:A:N6	2.31	0.45
1:CA:486:U:H2'	1:CA:487:A:H8	1.81	0.45
1:CA:51:A:H5''	1:CA:52:C:H5''	1.99	0.45
1:CA:613:C:H2'	1:CA:614:C:H6	1.82	0.45
1:CA:746:A:H2'	1:CA:747:A:C8	2.52	0.45
1:CA:818:G:H3'	1:CA:819:A:H5''	1.97	0.45
1:CA:932:C:H2'	1:CA:933:G:H8	1.81	0.45
1:CA:988:G:H21	1:CA:1015:G:H22	1.60	0.45
3:CD:148:ALA:C	3:CD:150:LYS:N	2.70	0.45
5:CF:7:VAL:HA	5:CF:60:VAL:O	2.17	0.45
7:CH:63:LYS:CG	7:CH:70:VAL:HG21	2.46	0.45
11:CL:82:ARG:NH2	11:CL:95:HIS:ND1	2.64	0.45
12:CM:33:LEU:HD13	12:CM:40:GLU:HA	1.97	0.45
21:CN:19:TYR:O	21:CN:21:ALA:N	2.48	0.45
13:CP:75:ILE:HG22	13:CP:80:LYS:HD2	1.98	0.45
15:CR:28:LEU:C	15:CR:30:ASN:H	2.19	0.45
15:CR:34:GLU:HB2	19:CU:18:PHE:CZ	2.52	0.45
16:CS:43:MET:O	16:CS:46:LEU:HB2	2.17	0.45
23:DB:1219:U:H2'	23:DB:1220:G:C8	2.52	0.45
23:DB:1526:C:O2'	23:DB:1527:G:H5'	2.17	0.45
23:DB:1593:A:H2'	23:DB:1594:U:H6	1.79	0.45
23:DB:2028:U:O2'	23:DB:2029:G:H5'	2.16	0.45
23:DB:2065:C:H2'	23:DB:2066:C:H6	1.78	0.45
23:DB:2512:C:P	26:DD:128:ARG:HB2	2.57	0.45
23:DB:2600:A:O2'	23:DB:2601:C:H5'	2.17	0.45
23:DB:592:A:H2'	23:DB:593:U:C6	2.50	0.45
23:DB:68:G:H2'	23:DB:69:C:H6	1.82	0.45
23:DB:909:A:H2'	23:DB:912:C:C5	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:921:C:H2'	23:DB:922:C:C6	2.52	0.45
23:DB:934:U:H2'	23:DB:935:C:H6	1.80	0.45
23:DB:969:G:H2'	23:DB:970:U:O4'	2.16	0.45
25:DC:110:LYS:O	25:DC:113:ASP:HB2	2.16	0.45
29:DE:137:LYS:O	29:DE:141:MET:HG3	2.16	0.45
47:DF:137:PHE:O	47:DF:138:PRO:C	2.55	0.45
47:DF:37:MET:SD	47:DF:52:ALA:HA	2.57	0.45
48:DG:169:ARG:HB3	48:DG:171:LYS:NZ	2.32	0.45
40:DH:141:LYS:O	40:DH:142:VAL:HG23	2.17	0.45
24:DI:23:VAL:HG12	24:DI:27:LEU:HD21	1.97	0.45
37:DL:122:VAL:HB	37:DL:143:GLU:OE2	2.16	0.45
43:DO:100:HIS:HA	43:DO:104:GLN:HG3	1.98	0.45
22:DA:7:G:H5''	43:DO:29:HIS:CD2	2.52	0.45
28:DP:50:ARG:HB3	28:DP:57:ALA:N	2.32	0.45
52:DW:23:LYS:CG	52:DW:24:ARG:N	2.78	0.45
1:AA:1101:A:O2'	1:AA:1102:A:OP2	2.34	0.45
1:AA:1265:C:O2'	1:AA:1266:G:H5'	2.17	0.45
1:AA:35:G:H2'	1:AA:36:C:H6	1.82	0.45
1:AA:437:U:C5'	3:AD:151:GLN:HE21	2.30	0.45
1:AA:605:U:H2'	1:AA:606:G:H8	1.82	0.45
1:AA:916:U:H2'	1:AA:917:G:H8	1.82	0.45
1:AA:975:A:H5''	1:AA:976:G:O5'	2.16	0.45
1:AA:1073:U:H4'	18:AB:104:LYS:CE	2.46	0.45
18:AB:15:PHE:O	18:AB:40:ILE:HD12	2.17	0.45
3:AD:187:ARG:O	3:AD:191:SER:N	2.50	0.45
3:AD:48:SER:O	3:AD:49:ASP:C	2.55	0.45
5:AF:34:GLY:O	5:AF:35:LYS:HB2	2.16	0.45
6:AG:14:ASP:CG	6:AG:17:PHE:HB2	2.36	0.45
8:AI:22:PRO:HA	8:AI:60:LEU:CB	2.46	0.45
8:AI:78:ILE:O	8:AI:82:ILE:HG13	2.17	0.45
8:AI:80:HIS:CE1	8:AI:84:ARG:HD3	2.51	0.45
8:AI:96:GLU:HA	8:AI:99:LYS:CE	2.47	0.45
34:B3:18:LYS:HD2	34:B3:19:GLY:H	1.80	0.45
34:B3:22:LYS:HA	34:B3:48:MET:HA	1.99	0.45
23:BB:1072:C:N3	23:BB:1092:C:N4	2.65	0.45
23:BB:1150:C:H2'	23:BB:1151:A:C8	2.50	0.45
23:BB:1564:C:H2'	23:BB:1565:C:C6	2.51	0.45
23:BB:1695:G:N7	25:BC:13:ARG:NH2	2.65	0.45
23:BB:1719:G:O2'	23:BB:1720:U:H5'	2.17	0.45
23:BB:1735:A:H2'	23:BB:1736:U:C6	2.52	0.45
23:BB:1784:A:H4'	23:BB:1785:A:O5'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1948:G:O2'	23:BB:1949:G:H5'	2.17	0.45
23:BB:2064:C:H1'	23:BB:2450:A:C6	2.52	0.45
23:BB:2655:G:H1'	23:BB:2656:U:H5	1.82	0.45
23:BB:2730:C:H2'	23:BB:2731:G:C8	2.50	0.45
23:BB:2847:U:H5''	28:BP:94:ALA:HB3	1.99	0.45
23:BB:2893:A:H4'	23:BB:2894:G:C5'	2.47	0.45
23:BB:406:G:O2'	23:BB:407:G:H5'	2.17	0.45
23:BB:742:A:O2'	23:BB:743:A:H5'	2.17	0.45
23:BB:842:U:H2'	23:BB:843:G:C8	2.49	0.45
25:BC:93:VAL:HG13	25:BC:101:ARG:H	1.81	0.45
26:BD:114:LYS:HD2	26:BD:116:LYS:CE	2.47	0.45
26:BD:8:LYS:HB2	26:BD:201:LEU:HD21	1.99	0.45
26:BD:51:THR:HG22	26:BD:52:THR:N	2.32	0.45
26:BD:55:LYS:HZ3	26:BD:56:LYS:CG	2.28	0.45
29:BE:48:THR:C	29:BE:50:ALA:N	2.70	0.45
47:BF:7:TYR:OH	47:BF:29:ARG:HG3	2.17	0.45
47:BF:34:THR:OG1	47:BF:154:THR:HB	2.16	0.45
48:BG:17:LYS:HE2	48:BG:19:ASN:CG	2.37	0.45
48:BG:15:ASP:CB	48:BG:26:LYS:H	2.29	0.45
27:BK:60:ALA:HA	27:BK:87:LEU:CD2	2.47	0.45
38:BM:105:MET:HB2	38:BM:117:PHE:CZ	2.48	0.45
38:BM:31:PHE:HA	38:BM:131:VAL:O	2.17	0.45
44:BQ:73:ILE:HG21	44:BQ:109:VAL:HG13	1.98	0.45
45:BS:73:LYS:HA	45:BS:73:LYS:HD2	1.68	0.45
50:BT:64:LYS:HE3	50:BT:64:LYS:N	2.32	0.45
35:BV:77:VAL:CA	35:BV:89:ILE:HG22	2.43	0.45
1:CA:1493:A:H2'	1:CA:1494:G:OP2	2.17	0.45
1:CA:420:U:H2'	1:CA:422:C:C4	2.52	0.45
1:CA:453:G:H2'	1:CA:454:G:C8	2.52	0.45
1:CA:583:A:N6	1:CA:758:C:H1'	2.32	0.45
18:CB:119:GLN:HA	18:CB:122:ASP:OD2	2.16	0.45
18:CB:93:HIS:CD2	18:CB:94:ARG:HH22	2.34	0.45
2:CC:129:PHE:O	2:CC:133:MET:HB3	2.17	0.45
3:CD:107:GLY:O	3:CD:157:ALA:HB1	2.17	0.45
5:CF:91:ARG:N	5:CF:93:LYS:HZ1	2.13	0.45
1:CA:1379:G:O6	6:CG:2:ARG:HD3	2.16	0.45
6:CG:63:VAL:C	6:CG:65:LEU:N	2.71	0.45
8:CI:117:LEU:N	8:CI:117:LEU:HD23	2.32	0.45
12:CM:13:HIS:N	12:CM:13:HIS:CD2	2.84	0.45
22:DA:60:C:H2'	22:DA:61:G:C8	2.45	0.45
23:DB:1359:A:H2'	23:DB:1360:G:O4'	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1409:U:H2'	23:DB:1410:G:H8	1.81	0.45
23:DB:1681:G:N2	23:DB:1762:A:H3'	2.32	0.45
23:DB:2210:U:C4	23:DB:2212:A:N7	2.85	0.45
23:DB:2064:C:H1'	23:DB:2450:A:C6	2.51	0.45
23:DB:2581:G:C8	23:DB:2610:C:N4	2.84	0.45
23:DB:2653:U:H3'	23:DB:2654:A:H2'	1.99	0.45
23:DB:2720:U:H2'	23:DB:2721:A:C8	2.51	0.45
23:DB:2808:G:O2'	23:DB:2809:A:H8	2.00	0.45
23:DB:2853:C:H2'	23:DB:2854:G:H8	1.82	0.45
23:DB:288:U:H2'	23:DB:289:G:C8	2.52	0.45
23:DB:43:G:H2'	23:DB:44:A:O4'	2.17	0.45
23:DB:664:G:H2'	23:DB:665:U:C6	2.50	0.45
25:DC:161:VAL:HG12	25:DC:173:LEU:HB2	1.99	0.45
23:DB:1799:G:OP1	25:DC:257:ARG:HG3	2.17	0.45
26:DD:38:LYS:HB2	26:DD:47:ALA:HB3	1.98	0.45
29:DE:111:GLU:HA	29:DE:114:ARG:CG	2.39	0.45
47:DF:116:LEU:CG	47:DF:117:SER:H	2.30	0.45
48:DG:15:ASP:CB	48:DG:26:LYS:H	2.29	0.45
40:DH:84:ALA:HB3	40:DH:148:ALA:CB	2.47	0.45
41:DJ:115:GLY:HA2	41:DJ:118:MET:HE3	1.99	0.45
41:DJ:73:VAL:HG23	41:DJ:74:TYR:H	1.81	0.45
37:DL:93:ASN:O	37:DL:95:LEU:HD12	2.17	0.45
38:DM:41:LEU:O	38:DM:93:VAL:HA	2.17	0.45
42:DN:38:LEU:CB	42:DN:39:PRO:HD3	2.45	0.45
28:DP:95:LYS:HG3	28:DP:97:TYR:HE1	1.82	0.45
44:DQ:63:ARG:HH21	44:DQ:64:ILE:CD1	2.26	0.45
49:DR:28:ALA:HB3	49:DR:31:GLU:CG	2.47	0.45
50:DT:53:VAL:CG1	50:DT:87:LEU:HB3	2.46	0.45
46:DU:66:VAL:C	46:DU:68:ASN:H	2.19	0.45
30:DY:50:VAL:HB	30:DY:53:MET:HB2	1.99	0.45
51:DZ:11:ARG:HB3	51:DZ:12:PRO:HD2	1.99	0.45
1:AA:983:A:H2	1:AA:1222:G:H22	1.65	0.45
1:AA:1365:G:H2'	1:AA:1366:C:C6	2.51	0.45
1:AA:32:A:H2'	1:AA:33:A:H8	1.79	0.45
1:AA:398:U:H2'	1:AA:399:G:C8	2.52	0.45
1:AA:976:G:OP1	21:AN:70:HIS:HA	2.17	0.45
2:AC:155:ARG:H	2:AC:162:ALA:CA	2.30	0.45
3:AD:138:PRO:HA	3:AD:181:PHE:HD2	1.82	0.45
3:AD:77:GLU:OE1	3:AD:80:ARG:HD3	2.17	0.45
7:AH:86:LYS:HG2	7:AH:124:ILE:HD11	1.98	0.45
10:AK:126:ARG:HA	10:AK:126:ARG:NE	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:149:ALA:HB2	10:AK:60:PHE:HB3	1.98	0.45
11:AL:24:GLU:C	11:AL:26:CYS:N	2.70	0.45
16:AS:20:LYS:HD2	16:AS:20:LYS:O	2.17	0.45
16:AS:38:THR:HG23	16:AS:69:LYS:HD3	1.99	0.45
31:B0:16:ARG:HA	31:B0:19:ASP:HB2	1.99	0.45
31:B0:51:ARG:HB2	31:B0:52:LYS:H	1.62	0.45
22:BA:93:C:H2'	22:BA:94:A:H8	1.82	0.45
23:BB:1275:A:H2'	23:BB:1276:A:O4'	2.17	0.45
23:BB:1718:G:H2'	23:BB:1719:G:H8	1.82	0.45
23:BB:1824:G:H2'	23:BB:1825:U:C6	2.52	0.45
23:BB:2018:G:O2'	23:BB:2019:A:H5'	2.17	0.45
23:BB:2145:C:H4'	23:BB:2145:C:OP2	2.17	0.45
23:BB:577:G:OP1	23:BB:2502:G:H2'	2.17	0.45
23:BB:1750:G:O2'	23:BB:2860:A:H2	2.00	0.45
23:BB:493:G:O2'	23:BB:494:G:H5'	2.17	0.45
23:BB:639:U:H2'	23:BB:640:C:H6	1.82	0.45
23:BB:783:A:H4'	23:BB:1779:U:O2	2.17	0.45
25:BC:109:LEU:H	25:BC:109:LEU:CD2	2.30	0.45
25:BC:124:LYS:HG3	25:BC:125:PRO:HD2	1.99	0.45
47:BF:137:PHE:O	47:BF:138:PRO:C	2.55	0.45
48:BG:16:VAL:HG13	48:BG:49:LEU:HD13	1.98	0.45
48:BG:5:LYS:HE3	48:BG:61:TRP:CZ3	2.51	0.45
41:BJ:101:ILE:O	41:BJ:105:VAL:HG22	2.17	0.45
41:BJ:19:ASP:CB	41:BJ:57:LEU:HB2	2.47	0.45
41:BJ:58:ASN:HA	41:BJ:127:GLY:HA2	1.99	0.45
37:BL:127:VAL:HG22	37:BL:131:ALA:HB3	1.99	0.45
37:BL:81:ASP:HA	37:BL:84:LYS:HE2	1.98	0.45
38:BM:41:LEU:O	38:BM:94:ALA:N	2.50	0.45
38:BM:64:TRP:C	38:BM:65:ILE:HG13	2.36	0.45
38:BM:65:ILE:HG23	38:BM:103:TYR:CD2	2.52	0.45
43:BO:37:ALA:HB1	43:BO:78:VAL:HG21	1.99	0.45
28:BP:98:TYR:C	28:BP:100:ARG:H	2.19	0.45
44:BQ:38:VAL:O	44:BQ:39:ILE:C	2.53	0.45
49:BR:3:ALA:HB2	49:BR:101:ILE:HD12	1.99	0.45
45:BS:3:THR:HB	45:BS:62:ASP:CB	2.47	0.45
1:CA:1008:U:H2'	1:CA:1009:U:C4'	2.47	0.45
1:CA:1430:A:H2'	1:CA:1431:A:O4'	2.17	0.45
1:CA:1492:A:H5''	1:CA:1493:A:N9	2.31	0.45
1:CA:259:G:O2'	1:CA:260:G:H5'	2.17	0.45
1:CA:371:A:O2'	1:CA:372:C:H5'	2.17	0.45
1:CA:551:U:H2'	1:CA:552:U:H6	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:69:G:N2	1:CA:71:A:H62	2.15	0.45
1:CA:895:G:H2'	1:CA:896:C:H6	1.82	0.45
1:CA:974:A:H4'	1:CA:975:A:H3'	1.98	0.45
2:CC:152:VAL:HG11	2:CC:156:LEU:HG	1.99	0.45
2:CC:67:ILE:O	2:CC:103:ALA:N	2.50	0.45
4:CE:81:GLN:N	4:CE:146:MET:HE3	2.18	0.45
7:CH:17:GLN:HE21	7:CH:62:LEU:CG	2.28	0.45
8:CI:14:SER:HA	8:CI:68:GLY:O	2.17	0.45
11:CL:35:ARG:HA	11:CL:35:ARG:CZ	2.47	0.45
21:CN:3:GLN:O	21:CN:6:LYS:HB2	2.17	0.45
13:CP:23:ASP:OD1	13:CP:25:ARG:HB2	2.16	0.45
23:DB:1032:A:OP1	32:D4:8:LYS:HG3	2.16	0.45
23:DB:1106:G:O2'	23:DB:1107:G:H5'	2.17	0.45
23:DB:1203:U:H3'	23:DB:1204:A:H5''	1.97	0.45
23:DB:1449:G:O2'	23:DB:1450:G:H5'	2.17	0.45
23:DB:2503:A:O2'	23:DB:2505:G:OP2	2.35	0.45
23:DB:2868:A:H2'	23:DB:2869:G:C8	2.52	0.45
23:DB:2896:C:H2'	23:DB:2897:U:C6	2.52	0.45
23:DB:302:C:H2'	23:DB:303:G:C8	2.50	0.45
23:DB:567:U:H2'	23:DB:568:U:O4'	2.17	0.45
23:DB:635:C:H2'	23:DB:636:G:C8	2.52	0.45
23:DB:657:U:H2'	23:DB:658:U:C6	2.52	0.45
23:DB:842:U:H2'	23:DB:843:G:C8	2.49	0.45
47:DF:128:SER:CB	47:DF:154:THR:HA	2.47	0.45
22:DA:43:C:H1'	47:DF:91:ARG:NH2	2.32	0.45
47:DF:32:LYS:N	47:DF:95:MET:SD	2.90	0.45
48:DG:156:TYR:O	48:DG:157:LYS:HB2	2.17	0.45
48:DG:17:LYS:O	48:DG:23:ILE:HG23	2.17	0.45
40:DH:53:GLU:O	40:DH:57:LYS:HB3	2.17	0.45
24:DI:131:THR:O	24:DI:135:MET:HG3	2.16	0.45
24:DI:138:VAL:HG12	24:DI:139:VAL:N	2.32	0.45
24:DI:17:ALA:C	24:DI:19:PRO:HD3	2.37	0.45
38:DM:71:LYS:CB	38:DM:93:VAL:HG12	2.46	0.45
28:DP:12:MET:HG2	28:DP:54:LEU:HA	1.98	0.45
44:DQ:115:ALA:C	44:DQ:117:ALA:H	2.20	0.45
44:DQ:96:ASP:C	44:DQ:98:ALA:H	2.19	0.45
23:DB:993:G:H1'	49:DR:91:GLN:NE2	2.31	0.45
45:DS:49:LYS:C	45:DS:51:LEU:H	2.19	0.45
50:DT:43:ILE:HG23	50:DT:58:VAL:HG21	1.99	0.45
35:DV:53:LYS:HZ2	35:DV:54:ALA:N	2.11	0.45
23:DB:2365:G:O2'	52:DW:59:PHE:CE1	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DX:52:ARG:O	39:DX:56:LEU:HD23	2.17	0.45
39:DX:7:ARG:O	39:DX:9:LYS:HD3	2.17	0.45
51:DZ:5:CYS:SG	51:DZ:8:THR:HG23	2.57	0.45
1:AA:1025:U:H6	1:AA:1025:U:O5'	2.00	0.44
1:AA:1300:G:O6	1:AA:1334:G:H3'	2.17	0.44
1:AA:1269:A:H2	1:AA:1312:G:N3	2.15	0.44
1:AA:1326:U:O2'	1:AA:1327:C:H5'	2.17	0.44
1:AA:261:U:H2'	1:AA:263:A:OP2	2.16	0.44
1:AA:300:A:H2'	1:AA:301:G:O4'	2.16	0.44
1:AA:411:A:H62	1:AA:413:G:H21	1.64	0.44
1:AA:551:U:H2'	1:AA:552:U:H6	1.81	0.44
1:AA:735:C:H2'	1:AA:736:C:C6	2.52	0.44
1:AA:780:A:O2'	1:AA:781:A:H5''	2.16	0.44
1:AA:852:G:H2'	1:AA:853:C:C6	2.52	0.44
18:AB:134:LEU:HD23	18:AB:134:LEU:N	2.33	0.44
2:AC:108:PRO:C	2:AC:110:LEU:H	2.20	0.44
2:AC:20:THR:HG23	2:AC:20:THR:O	2.17	0.44
3:AD:28:ASP:CB	3:AD:33:ILE:HG21	2.47	0.44
3:AD:61:ARG:HD2	3:AD:68:GLU:HA	1.99	0.44
5:AF:21:MET:HB3	5:AF:25:TYR:CZ	2.52	0.44
5:AF:7:VAL:HA	5:AF:60:VAL:O	2.17	0.44
5:AF:62:MET:HG3	5:AF:64:VAL:CG2	2.43	0.44
6:AG:49:LEU:HD12	6:AG:124:SER:OG	2.17	0.44
9:AJ:15:HIS:O	9:AJ:18:ILE:HG22	2.17	0.44
10:AK:16:SER:CA	10:AK:78:ILE:HA	2.46	0.44
13:AP:19:VAL:H	13:AP:38:PHE:HA	1.82	0.44
16:AS:14:LEU:CD1	16:AS:32:THR:HG23	2.47	0.44
19:AU:42:THR:O	19:AU:45:LYS:N	2.50	0.44
31:B0:41:HIS:HB2	42:BN:99:LYS:C	2.37	0.44
22:BA:52:A:H2'	22:BA:53:A:O4'	2.16	0.44
23:BB:1061:U:H5'	24:BI:9:LYS:HZ1	1.82	0.44
23:BB:1227:G:O2'	23:BB:1228:G:H5'	2.17	0.44
23:BB:136:G:H2'	23:BB:137:U:H1'	1.99	0.44
23:BB:1591:A:H2'	23:BB:1592:C:O4'	2.17	0.44
23:BB:170:U:O2'	23:BB:171:U:H5'	2.17	0.44
23:BB:1820:U:C4	25:BC:158:GLY:HA3	2.52	0.44
23:BB:195:A:H1'	23:BB:250:G:H21	1.82	0.44
23:BB:335:C:O2'	23:BB:336:C:H5'	2.17	0.44
23:BB:611:C:H2'	23:BB:612:G:O4'	2.17	0.44
23:BB:812:C:H2'	23:BB:813:U:H6	1.81	0.44
26:BD:106:LYS:CB	26:BD:206:ALA:H	2.24	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:17:THR:C	29:BE:19:PHE:H	2.20	0.44
29:BE:112:LEU:HD11	29:BE:180:LEU:HB3	1.99	0.44
47:BF:104:THR:C	47:BF:108:PRO:HG2	2.37	0.44
47:BF:109:ARG:HG3	47:BF:137:PHE:CB	2.43	0.44
47:BF:116:LEU:CG	47:BF:117:SER:H	2.30	0.44
48:BG:10:VAL:C	48:BG:14:VAL:HG21	2.37	0.44
24:BI:44:LYS:O	24:BI:48:ILE:HG13	2.16	0.44
41:BJ:58:ASN:CA	41:BJ:127:GLY:HA2	2.47	0.44
27:BK:18:ARG:HB2	27:BK:45:GLU:HG2	1.99	0.44
37:BL:109:LYS:HB2	37:BL:111:ILE:HD11	1.98	0.44
38:BM:113:ALA:O	38:BM:116:ALA:HB3	2.16	0.44
43:BO:111:ARG:HD2	43:BO:117:PHE:CD1	2.51	0.44
44:BQ:63:ARG:HH21	44:BQ:64:ILE:CD1	2.27	0.44
44:BQ:65:ASN:O	44:BQ:69:ARG:N	2.46	0.44
44:BQ:65:ASN:HA	44:BQ:75:TYR:HB2	1.99	0.44
45:BS:60:HIS:ND1	45:BS:60:HIS:O	2.50	0.44
52:BW:28:GLU:H	52:BW:31:LEU:HD12	1.82	0.44
52:BW:53:GLY:O	52:BW:56:HIS:N	2.46	0.44
51:BZ:11:ARG:HB3	51:BZ:12:PRO:HD2	1.99	0.44
51:BZ:53:ALA:C	51:BZ:55:GLY:H	2.21	0.44
1:CA:1052:U:H2'	1:CA:1055:A:OP1	2.17	0.44
1:CA:769:G:H4'	1:CA:1513:A:H4'	1.99	0.44
1:CA:430:A:OP2	3:CD:6:PRO:HA	2.17	0.44
1:CA:511:C:O2'	1:CA:512:U:H6	2.00	0.44
1:CA:520:A:N1	1:CA:536:C:H1'	2.32	0.44
1:CA:83:C:H1'	1:CA:84:U:C6	2.51	0.44
2:CC:148:ILE:HG13	2:CC:201:ILE:HG23	1.98	0.44
3:CD:61:ARG:HG3	3:CD:71:PHE:CD2	2.52	0.44
5:CF:16:GLU:H	5:CF:16:GLU:CD	2.20	0.44
6:CG:85:GLN:OE1	6:CG:85:GLN:HA	2.17	0.44
7:CH:48:PHE:HA	7:CH:59:GLU:O	2.17	0.44
32:D4:31:PRO:O	32:D4:34:LYS:HB3	2.17	0.44
22:DA:43:C:H5'	47:DF:62:GLN:HB3	1.98	0.44
23:DB:1291:C:O2'	23:DB:1292:G:H5'	2.17	0.44
23:DB:1292:G:H2'	23:DB:1293:C:C6	2.52	0.44
23:DB:1671:U:H2'	23:DB:1673:G:OP2	2.17	0.44
23:DB:1997:C:O2'	23:DB:1998:A:H5'	2.16	0.44
23:DB:2088:A:H2'	23:DB:2089:C:C6	2.51	0.44
23:DB:2221:G:H2'	23:DB:2222:C:C6	2.52	0.44
23:DB:2249:U:N3	23:DB:2253:G:OP2	2.43	0.44
23:DB:2485:G:O2'	23:DB:2486:C:H5'	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2693:G:H2'	23:DB:2694:G:C8	2.51	0.44
23:DB:2860:A:H2'	23:DB:2861:U:O4'	2.17	0.44
23:DB:39:G:O2'	23:DB:40:U:H5'	2.17	0.44
23:DB:992:C:O2'	23:DB:993:G:H5'	2.18	0.44
25:DC:173:LEU:H	25:DC:173:LEU:HD22	1.82	0.44
23:DB:1813:G:H1'	25:DC:49:THR:OG1	2.17	0.44
26:DD:118:PHE:O	26:DD:119:ALA:CB	2.62	0.44
26:DD:62:LYS:H	26:DD:62:LYS:HD2	1.82	0.44
41:DJ:46:PRO:C	41:DJ:48:VAL:H	2.20	0.44
42:DN:103:ARG:HB2	42:DN:110:MET:CE	2.45	0.44
44:DQ:105:PHE:CA	44:DQ:108:LEU:HD12	2.37	0.44
49:DR:72:VAL:CG2	49:DR:89:HIS:HB3	2.44	0.44
45:DS:17:VAL:C	45:DS:19:LEU:N	2.67	0.44
46:DU:64:ILE:HD11	46:DU:68:ASN:ND2	2.33	0.44
51:DZ:5:CYS:CB	51:DZ:10:LYS:H	2.30	0.44
1:AA:1017:U:H2'	1:AA:1018:G:C8	2.52	0.44
3:AD:167:PRO:HG2	3:AD:170:LEU:HD11	1.99	0.44
3:AD:53:GLN:HG2	3:AD:198:LEU:CD2	2.47	0.44
4:AE:131:ASN:O	4:AE:135:VAL:HG23	2.16	0.44
6:AG:74:VAL:HG12	6:AG:87:PRO:CD	2.47	0.44
13:AP:67:ILE:CG1	13:AP:72:ALA:HB2	2.46	0.44
14:AQ:80:LYS:NZ	14:AQ:80:LYS:H	2.15	0.44
17:AT:7:LYS:O	17:AT:11:ILE:HG13	2.17	0.44
33:B1:8:ILE:HG23	33:B1:9:LYS:N	2.32	0.44
34:B3:23:HIS:CG	37:BL:61:LEU:HD23	2.52	0.44
34:B3:30:HIS:HD2	34:B3:31:ILE:H	1.65	0.44
34:B3:41:ARG:C	34:B3:43:LEU:H	2.21	0.44
22:BA:102:G:H2'	22:BA:103:U:H6	1.82	0.44
23:BB:1745:A:O2'	23:BB:1746:A:H5'	2.17	0.44
23:BB:1877:A:H2'	23:BB:1878:G:H8	1.81	0.44
23:BB:2094:A:H4'	40:BH:25:TYR:CZ	2.53	0.44
23:BB:2357:G:N2	23:BB:2359:C:H3'	2.33	0.44
23:BB:2842:G:O2'	23:BB:2843:G:H5'	2.18	0.44
23:BB:2886:A:H2'	23:BB:2887:A:O4'	2.17	0.44
23:BB:57:C:H2'	23:BB:58:G:C8	2.52	0.44
23:BB:714:U:O2	23:BB:716:A:H3'	2.18	0.44
25:BC:102:TYR:C	25:BC:103:ILE:HG13	2.37	0.44
29:BE:58:LYS:NZ	29:BE:58:LYS:N	2.53	0.44
47:BF:126:ASN:HD22	47:BF:156:THR:HA	1.82	0.44
47:BF:121:PHE:CE1	47:BF:162:ASP:HB2	2.51	0.44
47:BF:168:LEU:C	47:BF:170:ALA:H	2.19	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BF:116:LEU:O	47:BF:177:ARG:HB2	2.16	0.44
37:BL:23:ILE:H	37:BL:23:ILE:CD1	2.15	0.44
38:BM:110:GLU:N	38:BM:110:GLU:CD	2.71	0.44
42:BN:38:LEU:CB	42:BN:39:PRO:HD3	2.45	0.44
28:BP:19:PHE:HB2	28:BP:50:ARG:NH1	2.31	0.44
44:BQ:68:ALA:HB1	44:BQ:73:ILE:HG23	1.99	0.44
46:BU:49:PRO:HA	46:BU:53:GLN:HG3	1.98	0.44
52:BW:67:LYS:O	52:BW:68:PHE:HB2	2.17	0.44
30:BY:40:THR:O	30:BY:43:ILE:HG23	2.17	0.44
30:BY:4:ILE:HD12	30:BY:58:GLU:HG3	2.00	0.44
1:CA:1118:U:OP1	8:CI:10:ARG:HD2	2.18	0.44
1:CA:1124:G:H5'	9:CJ:37:ARG:NE	2.31	0.44
1:CA:1148:U:O2'	1:CA:1149:C:H5'	2.17	0.44
1:CA:1219:A:H2'	1:CA:1220:G:C8	2.52	0.44
1:CA:1230:C:O2'	1:CA:1231:G:H5'	2.17	0.44
1:CA:1255:G:O2'	1:CA:1258:G:H1'	2.16	0.44
1:CA:1362:A:H4'	1:CA:1363:A:OP1	2.17	0.44
1:CA:343:U:O2'	1:CA:344:A:H8	2.00	0.44
1:CA:35:G:H2'	1:CA:36:C:C6	2.53	0.44
1:CA:394:G:O2'	1:CA:395:C:H5'	2.17	0.44
1:CA:709:U:H2'	1:CA:710:G:H8	1.81	0.44
1:CA:728:A:H2'	1:CA:729:A:C8	2.51	0.44
1:CA:735:C:H2'	1:CA:736:C:C6	2.52	0.44
18:CB:64:GLY:O	18:CB:88:GLN:HA	2.18	0.44
4:CE:89:THR:HB	4:CE:134:ASN:ND2	2.32	0.44
5:CF:92:THR:HG22	5:CF:93:LYS:N	2.32	0.44
7:CH:99:GLY:CA	7:CH:129:ALA:HA	2.46	0.44
8:CI:30:ASN:C	8:CI:32:ARG:H	2.20	0.44
8:CI:47:VAL:O	8:CI:50:PRO:HD2	2.17	0.44
10:CK:14:GLN:HA	10:CK:76:TYR:O	2.17	0.44
10:CK:17:ASP:HB3	10:CK:80:ASN:OD1	2.17	0.44
11:CL:14:LYS:HZ3	11:CL:16:ALA:CB	2.29	0.44
11:CL:28:GLN:HE21	11:CL:28:GLN:HB3	1.52	0.44
20:CO:24:SER:HB3	20:CO:27:VAL:CG2	2.48	0.44
13:CP:71:VAL:HG13	13:CP:72:ALA:H	1.82	0.44
15:CR:37:LYS:HG2	15:CR:37:LYS:H	1.63	0.44
16:CS:62:THR:HG22	16:CS:63:ASP:N	2.26	0.44
19:CU:42:THR:O	19:CU:45:LYS:N	2.50	0.44
31:D0:55:ALA:O	31:D0:56:LYS:C	2.55	0.44
23:DB:1031:G:H4'	32:D4:6:SER:HB3	1.98	0.44
22:DA:73:A:N3	22:DA:73:A:H2'	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:DA:76:G:H2'	22:DA:77:U:C6	2.53	0.44
23:DB:1005:C:H2'	23:DB:1006:C:H6	1.82	0.44
23:DB:1153:C:H2'	23:DB:1154:G:C8	2.52	0.44
23:DB:819:A:OP2	23:DB:1187:G:N2	2.49	0.44
23:DB:1392:A:C6	23:DB:1393:A:C6	3.06	0.44
23:DB:1739:A:H2'	23:DB:1740:G:C8	2.52	0.44
23:DB:182:A:H2'	23:DB:183:C:H6	1.82	0.44
23:DB:2056:G:N2	31:D0:1:ALA:HA	2.32	0.44
23:DB:2070:A:H2'	23:DB:2071:A:H8	1.75	0.44
23:DB:571:U:O2'	49:DR:80:ARG:NH2	2.50	0.44
23:DB:611:C:H2'	23:DB:612:G:O4'	2.17	0.44
23:DB:672:C:H2'	23:DB:673:C:H6	1.82	0.44
23:DB:779:U:O2'	23:DB:780:G:H5'	2.18	0.44
23:DB:81:G:O2'	23:DB:82:U:H5'	2.17	0.44
23:DB:834:G:O2'	23:DB:835:C:H5'	2.18	0.44
25:DC:211:ARG:C	25:DC:213:ARG:H	2.20	0.44
26:DD:54:ALA:N	26:DD:76:GLY:HA2	2.32	0.44
26:DD:55:LYS:C	26:DD:57:ALA:H	2.20	0.44
48:DG:34:ARG:HH11	48:DG:34:ARG:N	2.15	0.44
40:DH:64:ALA:O	40:DH:68:ARG:HG2	2.18	0.44
24:DI:100:ILE:HG23	24:DI:104:GLN:OE1	2.18	0.44
41:DJ:101:ILE:O	41:DJ:104:ALA:HB3	2.17	0.44
41:DJ:102:GLU:HA	41:DJ:105:VAL:HG22	1.98	0.44
42:DN:61:ALA:C	42:DN:63:ARG:N	2.67	0.44
28:DP:114:ASN:HA	28:DP:114:ASN:HD22	1.59	0.44
28:DP:62:LYS:HB3	28:DP:69:VAL:CG2	2.47	0.44
28:DP:89:GLY:N	28:DP:112:ARG:NH1	2.66	0.44
44:DQ:60:TRP:HB3	44:DQ:92:LYS:O	2.18	0.44
49:DR:1:MET:C	49:DR:2:TYR:HD1	2.21	0.44
49:DR:76:LYS:HB2	49:DR:85:LYS:HB2	1.99	0.44
50:DT:45:ALA:HA	50:DT:48:GLN:HG2	1.99	0.44
35:DV:77:VAL:CA	35:DV:89:ILE:HG22	2.42	0.44
52:DW:28:GLU:H	52:DW:31:LEU:HD12	1.81	0.44
52:DW:69:GLU:HG3	52:DW:80:SER:HG	1.82	0.44
30:DY:53:MET:HE2	30:DY:53:MET:HA	1.99	0.44
51:DZ:5:CYS:HB3	51:DZ:10:LYS:N	2.31	0.44
51:DZ:54:LYS:O	51:DZ:57:ARG:HB2	2.18	0.44
1:AA:1120:C:O2'	1:AA:1121:U:H5'	2.17	0.44
1:AA:1142:G:H3'	1:AA:1143:G:C8	2.53	0.44
1:AA:1151:A:HO2'	1:AA:1152:A:H8	1.62	0.44
1:AA:1293:C:H2'	1:AA:1294:G:O4'	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1326:U:H2'	1:AA:1327:C:C6	2.52	0.44
1:AA:1342:C:O2'	8:AI:125:GLN:HB3	2.18	0.44
1:AA:814:A:N7	1:AA:816:A:C4	2.85	0.44
18:AB:114:LYS:O	18:AB:117:GLU:N	2.50	0.44
2:AC:57:GLU:HB2	2:AC:64:ARG:HB2	1.98	0.44
3:AD:161:ALA:HA	3:AD:164:ARG:HB3	2.00	0.44
6:AG:144:ALA:C	6:AG:146:ALA:H	2.20	0.44
8:AI:107:ALA:O	8:AI:108:ARG:C	2.55	0.44
8:AI:18:VAL:HG11	8:AI:82:ILE:CA	2.44	0.44
11:AL:65:TYR:HB3	11:AL:95:HIS:CD2	2.52	0.44
12:AM:14:ALA:HB1	12:AM:33:LEU:HD11	1.99	0.44
20:AO:31:LEU:O	20:AO:34:ALA:HB3	2.17	0.44
16:AS:39:ILE:CD1	16:AS:65:MET:HB3	2.48	0.44
32:B4:8:LYS:HG2	32:B4:9:LYS:H	1.81	0.44
22:BA:113:C:H2'	22:BA:114:C:H6	1.81	0.44
23:BB:814:C:C2	23:BB:1194:A:C2	3.05	0.44
23:BB:1538:G:O2'	23:BB:1539:U:H5'	2.17	0.44
23:BB:1600:C:C2'	23:BB:1601:G:H5'	2.48	0.44
23:BB:2221:G:H2'	23:BB:2222:C:C6	2.53	0.44
23:BB:2240:U:O2'	23:BB:2241:A:H5'	2.17	0.44
23:BB:11:C:N4	23:BB:2629:U:H3	2.15	0.44
23:BB:2679:A:H2'	23:BB:2680:U:C6	2.52	0.44
23:BB:481:G:O5'	46:BU:43:LYS:HE2	2.17	0.44
23:BB:577:G:O2'	23:BB:578:G:H5'	2.18	0.44
23:BB:590:A:H2'	23:BB:591:U:H6	1.74	0.44
47:BF:111:ARG:O	47:BF:112:ASP:CB	2.62	0.44
22:BA:43:C:H1'	47:BF:91:ARG:NH2	2.32	0.44
40:BH:116:ARG:HH22	40:BH:133:GLN:N	2.16	0.44
40:BH:2:GLN:HB2	40:BH:39:ALA:HB2	1.99	0.44
41:BJ:6:ALA:HB3	41:BJ:45:THR:CG2	2.44	0.44
44:BQ:48:ASP:C	44:BQ:50:ARG:N	2.70	0.44
49:BR:55:ASP:CG	49:BR:56:GLY:H	2.21	0.44
50:BT:92:ASN:HD22	50:BT:92:ASN:HA	1.53	0.44
46:BU:40:LEU:HA	46:BU:60:LYS:O	2.17	0.44
51:BZ:33:LEU:H	51:BZ:52:SER:HB3	1.82	0.44
51:BZ:5:CYS:HB3	51:BZ:10:LYS:N	2.32	0.44
1:CA:1062:U:H2'	1:CA:1063:C:C6	2.52	0.44
1:CA:1133:G:H2'	1:CA:1134:G:O4'	2.17	0.44
1:CA:1225:A:OP1	12:CM:100:ARG:HA	2.17	0.44
1:CA:1270:G:H2'	1:CA:1271:A:H8	1.81	0.44
1:CA:1355:G:H2'	1:CA:1356:G:C8	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:261:U:H2'	1:CA:263:A:OP2	2.17	0.44
1:CA:450:G:N7	1:CA:481:G:O6	2.50	0.44
1:CA:71:A:N6	1:CA:100:G:N7	2.66	0.44
1:CA:570:G:H1'	1:CA:820:U:C4	2.52	0.44
1:CA:932:C:H2'	1:CA:933:G:C8	2.52	0.44
3:CD:111:ALA:O	3:CD:114:ARG:HB2	2.17	0.44
3:CD:61:ARG:HD2	3:CD:68:GLU:HA	1.99	0.44
4:CE:43:GLY:O	4:CE:72:ASN:HA	2.17	0.44
4:CE:45:VAL:H	4:CE:71:ILE:HG22	1.83	0.44
4:CE:45:VAL:O	4:CE:71:ILE:HG22	2.17	0.44
5:CF:6:ILE:HD11	5:CF:8:PHE:CD2	2.43	0.44
5:CF:74:LEU:HG	5:CF:78:PHE:CE1	2.52	0.44
6:CG:70:PRO:HB3	6:CG:102:TRP:HH2	1.80	0.44
6:CG:2:ARG:HH11	6:CG:2:ARG:CB	2.30	0.44
7:CH:118:ALA:HB3	7:CH:120:LEU:HD22	1.99	0.44
7:CH:46:GLU:N	7:CH:63:LYS:HE3	2.32	0.44
8:CI:29:ILE:HG12	8:CI:64:ILE:HD13	2.00	0.44
8:CI:8:THR:HG1	8:CI:80:HIS:CE1	2.35	0.44
9:CJ:8:ILE:HD13	9:CJ:76:ILE:HG23	1.99	0.44
11:CL:18:SER:OG	11:CL:19:ASN:N	2.51	0.44
21:CN:27:LYS:HA	21:CN:31:SER:HB3	1.98	0.44
56:CA:1910:HOH:O	21:CN:3:GLN:HB2	2.16	0.44
10:CK:125:LYS:O	19:CU:33:ARG:NH1	2.50	0.44
19:CU:34:ARG:NE	19:CU:34:ARG:C	2.70	0.44
31:D0:38:LEU:HD13	31:D0:41:HIS:NE2	2.33	0.44
23:DB:1723:G:N7	23:DB:1737:G:N2	2.58	0.44
23:DB:1752:C:O2'	23:DB:1753:G:H5'	2.18	0.44
23:DB:1796:U:O2'	23:DB:1797:G:H5'	2.18	0.44
23:DB:2454:G:O2'	23:DB:2455:G:H5'	2.18	0.44
23:DB:590:A:H2'	23:DB:591:U:H6	1.76	0.44
23:DB:734:A:C2	23:DB:735:A:H1'	2.52	0.44
23:DB:830:G:H4'	23:DB:831:G:OP2	2.16	0.44
23:DB:857:G:H2'	23:DB:858:G:H5'	1.96	0.44
25:DC:16:VAL:N	25:DC:203:VAL:HG11	2.25	0.44
25:DC:75:ALA:HA	25:DC:96:LYS:HZ2	1.82	0.44
26:DD:8:LYS:HG3	26:DD:9:VAL:N	2.31	0.44
47:DF:103:ILE:H	47:DF:106:ALA:HB3	1.81	0.44
48:DG:149:ALA:C	48:DG:151:ARG:H	2.21	0.44
48:DG:155:PRO:HA	48:DG:170:THR:HA	1.98	0.44
48:DG:84:LYS:HB2	48:DG:132:LEU:H	1.81	0.44
24:DI:102:ARG:HG3	24:DI:141:ASP:CB	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:DI:2:LYS:O	24:DI:3:LYS:HG3	2.16	0.44
41:DJ:55:ILE:CG2	41:DJ:123:LYS:HB2	2.44	0.44
41:DJ:30:THR:N	41:DJ:108:MET:HE3	2.32	0.44
27:DK:58:LEU:HD11	27:DK:86:LEU:HB2	1.98	0.44
37:DL:121:THR:CB	37:DL:141:LYS:HB3	2.46	0.44
37:DL:68:SER:CB	37:DL:71:ALA:HB3	2.47	0.44
23:DB:1252:G:N2	44:DQ:32:ARG:HB3	2.32	0.44
44:DQ:67:ALA:HB1	44:DQ:105:PHE:CZ	2.52	0.44
44:DQ:79:ILE:C	44:DQ:79:ILE:HD13	2.38	0.44
44:DQ:96:ASP:C	44:DQ:98:ALA:N	2.70	0.44
50:DT:25:GLU:OE1	50:DT:30:ILE:HA	2.17	0.44
46:DU:73:ASN:ND2	46:DU:74:ALA:N	2.65	0.44
39:DX:6:LEU:C	39:DX:8:GLU:H	2.21	0.44
1:AA:1354:U:H6	1:AA:1354:U:O5'	2.00	0.44
1:AA:153:C:H2'	1:AA:154:U:C6	2.53	0.44
1:AA:764:C:O2'	1:AA:765:G:H5'	2.18	0.44
1:AA:978:A:H5'	1:AA:1224:U:O4	2.17	0.44
18:AB:107:ARG:HA	18:AB:110:ILE:CD1	2.47	0.44
8:AI:6:TYR:HA	8:AI:18:VAL:O	2.17	0.44
31:B0:42:ILE:HG22	31:B0:43:THR:O	2.18	0.44
23:BB:1151:A:H2'	23:BB:1152:C:O4'	2.18	0.44
23:BB:123:G:H2'	23:BB:124:G:C8	2.52	0.44
23:BB:1664:A:H1'	23:BB:2726:A:C2	2.53	0.44
23:BB:2346:A:O4'	23:BB:2383:G:O4'	2.36	0.44
23:BB:2581:G:N3	23:BB:2581:G:H2'	2.33	0.44
23:BB:2710:C:H2'	23:BB:2711:A:H8	1.83	0.44
23:BB:2848:G:H22	23:BB:2867:G:N2	2.15	0.44
23:BB:296:U:H2'	23:BB:297:G:C8	2.52	0.44
23:BB:322:A:H1'	23:BB:339:U:O2	2.16	0.44
23:BB:921:C:H2'	23:BB:922:C:C6	2.52	0.44
29:BE:172:ALA:HB3	29:BE:195:GLN:NE2	2.33	0.44
47:BF:128:SER:CB	47:BF:154:THR:HA	2.47	0.44
48:BG:84:LYS:O	48:BG:85:LYS:O	2.36	0.44
40:BH:124:THR:HG22	40:BH:125:THR:N	2.33	0.44
37:BL:116:VAL:HG13	37:BL:117:THR:N	2.18	0.44
42:BN:70:THR:HB	42:BN:75:ILE:HD11	1.99	0.44
43:BO:115:LEU:H	43:BO:115:LEU:HD22	1.82	0.44
46:BU:86:PHE:CE1	46:BU:90:LYS:HB2	2.52	0.44
22:BA:94:A:OP1	35:BV:19:ARG:HD3	2.16	0.44
52:BW:23:LYS:NZ	52:BW:24:ARG:HG3	2.32	0.44
52:BW:46:ALA:HB3	52:BW:79:ILE:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1009:U:H2'	1:CA:1010:U:C6	2.52	0.44
1:CA:1247:U:O2'	1:CA:1248:A:H5'	2.18	0.44
1:CA:1343:G:H2'	1:CA:1344:C:O4'	2.17	0.44
1:CA:1492:A:H5''	1:CA:1493:A:C5	2.53	0.44
1:CA:1519:A:H3'	1:CA:1520:C:C5'	2.47	0.44
1:CA:515:G:H2'	1:CA:516:U:O4'	2.18	0.44
1:CA:807:A:H2'	1:CA:808:C:C6	2.53	0.44
1:CA:852:G:H2'	1:CA:853:C:C6	2.53	0.44
1:CA:963:G:H2'	1:CA:964:A:H8	1.83	0.44
18:CB:150:ILE:O	18:CB:150:ILE:HG12	2.17	0.44
3:CD:117:VAL:HA	3:CD:122:ILE:HG13	1.98	0.44
6:CG:121:ASN:ND2	6:CG:121:ASN:N	2.66	0.44
7:CH:14:ARG:HG3	7:CH:15:ASN:N	2.33	0.44
21:CN:44:VAL:C	21:CN:46:LYS:H	2.20	0.44
21:CN:63:CYS:SG	21:CN:82:LYS:HG3	2.57	0.44
34:D3:16:THR:C	34:D3:18:LYS:H	2.21	0.44
22:DA:29:A:H3'	22:DA:30:C:H6	1.82	0.44
23:DB:1141:U:OP2	41:DJ:65:THR:HG21	2.17	0.44
23:DB:142:A:C2	50:DT:2:ILE:CG2	3.00	0.44
23:DB:2454:G:H2'	23:DB:2455:G:H5'	1.98	0.44
23:DB:2798:U:H4'	23:DB:2800:A:C2	2.53	0.44
23:DB:2839:G:O2'	23:DB:2840:C:H5'	2.18	0.44
23:DB:305:C:O2'	23:DB:306:U:H5'	2.18	0.44
23:DB:338:G:N2	23:DB:339:U:H1'	2.32	0.44
23:DB:418:C:H2'	23:DB:419:U:H6	1.83	0.44
23:DB:53:A:H2'	23:DB:54:G:O4'	2.17	0.44
23:DB:554:U:H2'	23:DB:555:G:O4'	2.18	0.44
23:DB:817:C:H2'	23:DB:818:G:C8	2.53	0.44
23:DB:917:A:H2'	23:DB:918:A:O4'	2.18	0.44
25:DC:61:TYR:HA	25:DC:85:ASN:HD21	1.82	0.44
29:DE:108:ILE:HD13	29:DE:108:ILE:O	2.16	0.44
47:DF:131:VAL:O	47:DF:133:GLU:N	2.50	0.44
48:DG:17:LYS:HB3	48:DG:24:THR:O	2.18	0.44
40:DH:48:GLU:CB	40:DH:51:ARG:HH21	2.30	0.44
41:DJ:38:GLY:O	41:DJ:43:GLU:HB2	2.18	0.44
37:DL:132:ARG:O	37:DL:136:GLU:HB2	2.17	0.44
34:D3:13:PHE:CZ	37:DL:61:LEU:HD22	2.53	0.44
38:DM:71:LYS:CG	38:DM:93:VAL:HG12	2.46	0.44
44:DQ:65:ASN:O	44:DQ:69:ARG:HB2	2.18	0.44
1:AA:1117:A:H2	1:AA:1180:A:H1'	1.83	0.44
1:AA:1307:U:H2'	1:AA:1308:U:C6	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:286:C:H2'	1:AA:287:U:C6	2.52	0.44
1:AA:435:A:N3	1:AA:435:A:H2'	2.32	0.44
1:AA:510:A:N3	1:AA:543:U:H1'	2.32	0.44
18:AB:67:LEU:O	18:AB:160:LEU:HD23	2.18	0.44
3:AD:21:LYS:C	3:AD:23:GLY:H	2.20	0.44
8:AI:27:ILE:CD1	8:AI:34:LEU:HD22	2.48	0.44
8:AI:84:ARG:O	8:AI:87:MET:HB3	2.17	0.44
12:AM:2:ARG:O	12:AM:3:ILE:HB	2.18	0.44
21:AN:61:ASN:O	21:AN:62:ARG:O	2.36	0.44
20:AO:81:LEU:O	20:AO:85:LEU:HD13	2.16	0.44
14:AQ:8:GLN:HB3	14:AQ:59:GLU:HG3	1.99	0.44
16:AS:6:LYS:HB2	21:AN:40:ARG:HH12	1.83	0.44
17:AT:63:LYS:HD3	17:AT:63:LYS:HA	1.85	0.44
19:AU:40:PRO:C	19:AU:42:THR:N	2.71	0.44
23:BB:2348:U:H4'	33:B1:40:PRO:HG2	2.00	0.44
23:BB:1080:A:O2'	24:BI:126:ARG:HB2	2.16	0.44
23:BB:1158:C:O3'	30:BY:30:ARG:NH2	2.51	0.44
23:BB:1249:U:O4'	44:BQ:3:VAL:HG21	2.18	0.44
23:BB:1474:U:H2'	23:BB:1475:G:H5'	1.99	0.44
23:BB:1681:G:N2	23:BB:1762:A:H3'	2.32	0.44
23:BB:1808:A:H3'	23:BB:1809:A:H8	1.81	0.44
23:BB:1920:C:H2'	23:BB:1921:G:C8	2.52	0.44
23:BB:2154:A:C2'	23:BB:2155:U:H5'	2.47	0.44
23:BB:2666:C:O4'	23:BB:2666:C:O2	2.34	0.44
23:BB:2740:A:H2'	23:BB:2741:A:C8	2.52	0.44
23:BB:2798:U:H4'	23:BB:2800:A:C2	2.52	0.44
23:BB:329:G:O6	46:BU:16:LYS:HB2	2.18	0.44
25:BC:149:LYS:HG2	25:BC:152:GLN:NE2	2.32	0.44
47:BF:2:LYS:NZ	47:BF:97:GLU:HA	2.33	0.44
48:BG:84:LYS:HB2	48:BG:132:LEU:H	1.83	0.44
48:BG:39:ALA:C	48:BG:54:ARG:HG3	2.38	0.44
40:BH:72:ILE:O	40:BH:72:ILE:HG23	2.15	0.44
24:BI:79:LEU:HD11	24:BI:131:THR:OG1	2.16	0.44
28:BP:19:PHE:CZ	28:BP:25:VAL:HG11	2.52	0.44
46:BU:13:LEU:CA	46:BU:18:LYS:HE3	2.46	0.44
46:BU:46:LYS:HE3	46:BU:47:PRO:O	2.17	0.44
52:BW:43:LYS:HE3	52:BW:68:PHE:HE1	1.82	0.44
1:CA:137:U:H2'	1:CA:138:G:C8	2.51	0.44
1:CA:1411:C:H2'	1:CA:1412:C:C6	2.53	0.44
1:CA:605:U:H2'	1:CA:606:G:H8	1.82	0.44
1:CA:656:G:H2'	1:CA:657:U:C6	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:750:C:O2	20:CO:23:GLY:HA3	2.17	0.44
18:CB:59:ILE:HG13	18:CB:59:ILE:H	1.53	0.44
2:CC:112:ALA:CA	2:CC:201:ILE:HD11	2.48	0.44
3:CD:12:ARG:HG2	3:CD:12:ARG:H	1.52	0.44
3:CD:161:ALA:HA	3:CD:164:ARG:HB3	1.99	0.44
5:CF:38:ARG:HE	5:CF:63:ASN:CG	2.21	0.44
6:CG:121:ASN:HD22	6:CG:121:ASN:N	2.15	0.44
6:CG:148:LYS:O	6:CG:151:ALA:N	2.50	0.44
7:CH:44:PHE:CE2	7:CH:100:ILE:HG12	2.49	0.44
10:CK:126:ARG:NE	10:CK:126:ARG:HA	2.32	0.44
11:CL:80:LEU:HB3	11:CL:97:VAL:HG23	1.99	0.44
12:CM:65:GLU:HB2	12:CM:66:GLY:H	1.58	0.44
1:CA:1317:C:OP1	21:CN:56:PRO:HD2	2.17	0.44
13:CP:72:ALA:HA	13:CP:75:ILE:HD12	2.00	0.44
16:CS:29:PRO:HA	16:CS:47:THR:CB	2.47	0.44
34:D3:22:LYS:HA	34:D3:48:MET:HA	1.99	0.44
22:DA:19:C:H2'	22:DA:20:G:H8	1.82	0.44
23:DB:111:A:H2'	23:DB:112:U:O4'	2.18	0.44
23:DB:1241:A:H3'	23:DB:1242:U:H6	1.82	0.44
23:DB:1295:C:H2'	23:DB:1296:G:H8	1.83	0.44
23:DB:1591:A:H2'	23:DB:1592:C:O4'	2.17	0.44
23:DB:1595:C:O2'	23:DB:1596:A:H5'	2.18	0.44
23:DB:1718:G:H2'	23:DB:1719:G:H8	1.82	0.44
23:DB:2331:G:O2'	23:DB:2332:C:H5'	2.18	0.44
23:DB:27:G:HO2'	23:DB:28:A:H8	1.62	0.44
23:DB:535:G:O2'	23:DB:536:G:H5'	2.18	0.44
23:DB:596:U:H2'	23:DB:597:G:C8	2.52	0.44
23:DB:711:G:O2'	23:DB:712:G:H5'	2.17	0.44
23:DB:1792:G:OP1	25:DC:204:LEU:HD13	2.17	0.44
25:DC:251:THR:O	25:DC:252:LYS:HD2	2.18	0.44
26:DD:11:MET:HE1	26:DD:192:ALA:H	1.83	0.44
48:DG:10:VAL:C	48:DG:14:VAL:HG21	2.37	0.44
40:DH:5:LEU:C	40:DH:7:ASP:H	2.21	0.44
40:DH:74:ALA:C	40:DH:76:GLU:H	2.20	0.44
27:DK:88:ASN:ND2	27:DK:89:ASN:N	2.62	0.44
38:DM:43:ALA:HB2	38:DM:92:TRP:O	2.18	0.44
44:DQ:9:ALA:C	44:DQ:11:ALA:N	2.70	0.44
49:DR:14:VAL:CG2	49:DR:18:GLN:HG3	2.47	0.44
23:DB:1161:C:H4'	49:DR:8:GLY:O	2.17	0.44
45:DS:41:LYS:O	45:DS:43:ALA:N	2.50	0.44
35:DV:38:LEU:CG	35:DV:40:ILE:HG23	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DW:32:ALA:C	52:DW:34:SER:N	2.69	0.44
51:DZ:5:CYS:CB	51:DZ:10:LYS:N	2.81	0.44
1:AA:1360:A:C2'	1:AA:1361:G:H5'	2.47	0.44
1:AA:1441:A:C2	28:BP:113:LEU:HD13	2.53	0.44
1:AA:465:A:O2'	1:AA:466:A:H2'	2.17	0.44
1:AA:583:A:N6	1:AA:758:C:H1'	2.33	0.44
3:AD:148:ALA:C	3:AD:150:LYS:N	2.71	0.44
4:AE:113:VAL:O	4:AE:114:LEU:C	2.55	0.44
8:AI:24:ASN:O	8:AI:60:LEU:N	2.51	0.44
10:AK:63:GLN:HG3	10:AK:98:ALA:HB1	2.00	0.44
21:AN:20:PHE:HA	21:AN:23:ARG:HB2	2.00	0.44
2:AC:28:PHE:HD1	21:AN:75:LYS:HE3	1.82	0.44
20:AO:39:LEU:HD22	20:AO:56:LEU:HD13	2.00	0.44
20:AO:29:VAL:HG13	20:AO:67:LEU:HD21	2.00	0.44
15:AR:34:GLU:N	15:AR:34:GLU:OE1	2.48	0.44
16:AS:25:GLY:O	16:AS:27:LYS:HG2	2.17	0.44
17:AT:64:GLY:O	17:AT:65:LEU:HB3	2.17	0.44
33:B1:36:LYS:HG3	33:B1:46:VAL:O	2.18	0.44
34:B3:40:LYS:O	34:B3:43:LEU:N	2.48	0.44
22:BA:67:G:O2'	22:BA:68:C:H5'	2.16	0.44
22:BA:76:G:H2'	22:BA:77:U:C6	2.51	0.44
23:BB:1076:C:H2'	23:BB:1077:A:C8	2.52	0.44
23:BB:132:G:H2'	23:BB:133:U:C6	2.52	0.44
23:BB:1418:G:H1'	23:BB:1580:A:H61	1.82	0.44
23:BB:1716:U:H2'	23:BB:1717:A:C8	2.53	0.44
23:BB:2197:U:O2'	23:BB:2198:A:H2'	2.17	0.44
23:BB:2361:G:H2'	23:BB:2362:C:H6	1.82	0.44
23:BB:2494:G:H2'	23:BB:2495:G:H8	1.82	0.44
23:BB:2543:G:H4'	23:BB:2645:G:N2	2.32	0.44
23:BB:2559:C:H2'	23:BB:2560:A:H8	1.82	0.44
23:BB:2635:A:H4'	26:BD:79:LEU:HB2	1.99	0.44
23:BB:2688:G:H1'	23:BB:2721:A:H61	1.82	0.44
23:BB:274:C:H2'	23:BB:275:C:H6	1.82	0.44
23:BB:2811:G:O2'	23:BB:2812:G:H5'	2.18	0.44
23:BB:70:G:H3'	23:BB:113:U:C4'	2.44	0.44
23:BB:711:G:O2'	23:BB:712:G:H5'	2.18	0.44
26:BD:183:GLU:N	26:BD:183:GLU:CD	2.71	0.44
26:BD:32:ASN:HA	26:BD:51:THR:O	2.18	0.44
23:BB:600:G:C1'	29:BE:100:MET:HG2	2.45	0.44
29:BE:148:ILE:HD13	29:BE:187:VAL:CG2	2.47	0.44
47:BF:103:ILE:H	47:BF:106:ALA:HB3	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:28:ASN:HA	40:BH:28:ASN:HD22	1.60	0.44
40:BH:84:ALA:CA	40:BH:90:LEU:HG	2.47	0.44
24:BI:4:VAL:HG13	24:BI:4:VAL:O	2.17	0.44
24:BI:5:GLN:HG2	24:BI:6:ALA:H	1.81	0.44
37:BL:134:ALA:O	37:BL:137:ALA:HB3	2.17	0.44
37:BL:81:ASP:HA	37:BL:84:LYS:CE	2.47	0.44
38:BM:40:ARG:NH2	38:BM:73:ILE:HD12	2.33	0.44
38:BM:66:ARG:HE	38:BM:101:VAL:HG21	1.83	0.44
42:BN:36:THR:OG1	42:BN:37:THR:N	2.50	0.44
42:BN:55:ALA:HA	42:BN:80:PHE:CE1	2.53	0.44
42:BN:79:LEU:HD23	42:BN:83:LEU:HB3	1.98	0.44
45:BS:20:VAL:HG13	45:BS:21:ALA:N	2.33	0.44
46:BU:3:LYS:O	46:BU:4:ILE:HD13	2.17	0.44
51:BZ:14:THR:HA	51:BZ:28:ARG:HB2	1.99	0.44
51:BZ:55:GLY:O	51:BZ:58:VAL:HB	2.18	0.44
1:CA:1038:C:O2'	1:CA:1039:G:H5'	2.18	0.44
1:CA:1496:C:H2'	1:CA:1497:G:O4'	2.18	0.44
1:CA:635:A:H2'	1:CA:636:U:H6	1.83	0.44
1:CA:672:U:H2'	1:CA:673:A:C8	2.52	0.44
1:CA:815:A:H4'	1:CA:817:C:C5	2.52	0.44
3:CD:61:ARG:NH1	3:CD:68:GLU:HB2	2.33	0.44
4:CE:113:VAL:O	4:CE:116:VAL:HG22	2.16	0.44
6:CG:50:ALA:HB2	6:CG:57:GLU:N	2.32	0.44
9:CJ:36:VAL:HG22	9:CJ:76:ILE:HB	2.00	0.44
11:CL:23:LEU:O	11:CL:25:ALA:N	2.51	0.44
11:CL:36:VAL:O	11:CL:36:VAL:HG23	2.18	0.44
21:CN:60:ARG:NE	21:CN:62:ARG:NE	2.66	0.44
21:CN:73:LEU:O	21:CN:74:ARG:C	2.56	0.44
14:CQ:4:ILE:O	14:CQ:4:ILE:HD12	2.17	0.44
15:CR:70:THR:HB	15:CR:72:ARG:NH1	2.32	0.44
23:DB:1222:U:P	49:DR:90:ARG:HH22	2.40	0.44
23:DB:125:A:C2	36:D2:10:LEU:HA	2.53	0.44
23:DB:1459:G:H3'	23:DB:1460:U:H4'	1.99	0.44
23:DB:1549:A:H2'	23:DB:1550:C:H6	1.81	0.44
23:DB:1716:U:H2'	23:DB:1717:A:C8	2.52	0.44
23:DB:1948:G:O2'	23:DB:1949:G:H5'	2.18	0.44
1:CA:1483:A:H2	23:DB:1959:G:HO2'	1.62	0.44
23:DB:202:U:H2'	23:DB:203:A:O4'	2.17	0.44
23:DB:2267:A:OP2	23:DB:2268:A:H5''	2.18	0.44
23:DB:2863:C:O2'	23:DB:2864:G:H5'	2.17	0.44
23:DB:2834:G:H1'	23:DB:2883:A:H61	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:829:A:H5'	23:DB:831:G:N7	2.32	0.44
23:DB:863:A:O2'	23:DB:864:G:H5'	2.18	0.44
23:DB:898:C:OP1	23:DB:899:A:H4'	2.16	0.44
23:DB:968:C:H2'	23:DB:969:G:C8	2.49	0.44
25:DC:102:TYR:C	25:DC:103:ILE:HG13	2.37	0.44
23:DB:2637:U:OP1	26:DD:83:ARG:HD3	2.18	0.44
47:DF:24:VAL:C	47:DF:26:GLN:H	2.21	0.44
47:DF:2:LYS:NZ	47:DF:97:GLU:HA	2.33	0.44
47:DF:73:VAL:O	47:DF:74:ALA:HB2	2.17	0.44
48:DG:16:VAL:HG13	48:DG:49:LEU:HD13	1.99	0.44
48:DG:71:LEU:HA	48:DG:74:MET:SD	2.58	0.44
40:DH:60:GLU:C	40:DH:62:LEU:H	2.20	0.44
24:DI:89:SER:HA	24:DI:97:VAL:HG11	2.00	0.44
41:DJ:56:VAL:HG21	41:DJ:101:ILE:HG21	2.00	0.44
38:DM:40:ARG:NH2	38:DM:73:ILE:HD12	2.33	0.44
43:DO:51:ALA:HB3	43:DO:78:VAL:HG13	2.00	0.44
44:DQ:50:ARG:N	44:DQ:50:ARG:HD2	2.33	0.44
45:DS:20:VAL:HG13	45:DS:21:ALA:N	2.32	0.44
31:D0:27:LEU:HD21	45:DS:38:TYR:CE1	2.52	0.44
50:DT:32:LEU:N	50:DT:83:ALA:CB	2.81	0.44
46:DU:49:PRO:HA	46:DU:53:GLN:HG3	1.99	0.44
35:DV:16:ALA:HA	35:DV:19:ARG:CZ	2.48	0.44
30:DY:16:LEU:O	30:DY:19:HIS:HB2	2.17	0.44
1:AA:1173:U:H2'	1:AA:1174:G:O4'	2.18	0.44
1:AA:1444:U:H2'	1:AA:1445:U:C6	2.52	0.44
1:AA:515:G:H2'	1:AA:516:U:O4'	2.17	0.44
1:AA:635:A:H2'	1:AA:636:U:H6	1.82	0.44
18:AB:80:LYS:HB3	18:AB:90:PHE:CE1	2.53	0.44
2:AC:173:PRO:HB2	2:AC:176:THR:OG1	2.18	0.44
2:AC:67:ILE:O	2:AC:103:ALA:N	2.48	0.44
3:AD:10:LEU:HD22	3:AD:62:ARG:NH1	2.33	0.44
3:AD:116:LEU:HD21	3:AD:153:ARG:HD3	1.99	0.44
5:AF:15:SER:HA	5:AF:18:VAL:HG23	1.99	0.44
1:AA:586:C:H5''	7:AH:81:GLY:HA2	2.00	0.44
8:AI:41:GLU:C	8:AI:43:ALA:H	2.21	0.44
11:AL:23:LEU:O	11:AL:25:ALA:N	2.50	0.44
2:AC:11:LEU:HB2	21:AN:96:LYS:HB2	1.99	0.44
13:AP:38:PHE:CE2	13:AP:51:ARG:HD3	2.53	0.44
17:AT:49:ALA:C	17:AT:52:GLU:HB3	2.38	0.44
36:B2:10:LEU:O	36:B2:14:ARG:HG2	2.18	0.44
22:BA:43:C:H5'	47:BF:62:GLN:HB3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:100:U:H1'	23:BB:101:A:C4	2.52	0.44
23:BB:1183:U:H2'	23:BB:1184:U:H6	1.82	0.44
23:BB:137:U:OP2	23:BB:137:U:H6	2.00	0.44
23:BB:1410:G:H2'	23:BB:1411:U:C6	2.52	0.44
23:BB:1792:G:O2'	23:BB:1793:C:H5'	2.18	0.44
23:BB:2056:G:N2	31:B0:1:ALA:HA	2.32	0.44
23:BB:19:A:H2'	23:BB:20:C:H6	1.83	0.44
23:BB:337:C:OP1	46:BU:3:LYS:HG3	2.18	0.44
23:BB:554:U:H2'	23:BB:555:G:O4'	2.18	0.44
23:BB:751:A:OP2	45:BS:90:LYS:N	2.48	0.44
23:BB:830:G:H4'	23:BB:831:G:OP2	2.17	0.44
25:BC:109:LEU:N	25:BC:109:LEU:HD23	2.33	0.44
25:BC:230:PRO:O	25:BC:241:LYS:HD3	2.18	0.44
25:BC:40:GLY:O	25:BC:53:ILE:HG23	2.18	0.44
25:BC:76:VAL:HG12	25:BC:114:GLN:HB2	2.00	0.44
26:BD:173:GLN:O	26:BD:174:SER:HB3	2.17	0.44
47:BF:107:VAL:HA	47:BF:111:ARG:HH12	1.82	0.44
47:BF:134:GLN:C	47:BF:136:ILE:N	2.71	0.44
40:BH:81:ALA:CA	40:BH:146:VAL:HA	2.46	0.44
40:BH:147:VAL:HG12	40:BH:148:ALA:N	2.29	0.44
41:BJ:56:VAL:HG21	41:BJ:101:ILE:HG21	2.00	0.44
41:BJ:46:PRO:C	41:BJ:48:VAL:H	2.21	0.44
27:BK:71:ARG:HB3	27:BK:72:PRO:CD	2.41	0.44
37:BL:111:ILE:HA	37:BL:128:THR:HG23	1.98	0.44
38:BM:71:LYS:CB	38:BM:93:VAL:HG12	2.47	0.44
28:BP:45:VAL:N	28:BP:60:VAL:HG13	2.32	0.44
44:BQ:104:ALA:HA	49:BR:46:GLU:OE1	2.17	0.44
49:BR:83:TYR:HE2	49:BR:85:LYS:HE3	1.82	0.44
45:BS:50:VAL:HG12	45:BS:105:VAL:HG23	2.00	0.44
45:BS:61:ASN:HB3	45:BS:62:ASP:H	1.63	0.44
45:BS:71:VAL:HG22	45:BS:71:VAL:O	2.17	0.44
46:BU:51:LEU:N	46:BU:53:GLN:NE2	2.65	0.44
30:BY:50:VAL:HG23	30:BY:54:VAL:HG22	2.00	0.44
1:CA:545:C:O2'	1:CA:546:A:H5'	2.18	0.44
1:CA:551:U:H2'	1:CA:552:U:C6	2.52	0.44
1:CA:754:C:H3'	1:CA:754:C:O2	2.18	0.44
18:CB:100:LEU:HD23	18:CB:174:GLU:O	2.18	0.44
18:CB:64:GLY:H	18:CB:224:ARG:HH11	1.66	0.44
18:CB:69:VAL:HG23	18:CB:162:VAL:HB	2.00	0.44
2:CC:149:LYS:HD2	2:CC:166:TRP:CZ2	2.53	0.44
3:CD:106:PHE:CD1	3:CD:144:ILE:HD11	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:157:ALA:C	3:CD:159:GLU:H	2.20	0.44
3:CD:18:LEU:HD22	3:CD:18:LEU:N	2.33	0.44
3:CD:26:ALA:HA	3:CD:30:LYS:CE	2.44	0.44
4:CE:93:VAL:HG12	4:CE:94:PHE:N	2.33	0.44
5:CF:34:GLY:O	5:CF:35:LYS:HB2	2.18	0.44
6:CG:130:LYS:HD3	6:CG:130:LYS:C	2.38	0.44
10:CK:60:PHE:CE1	6:CG:146:ALA:HB1	2.52	0.44
9:CJ:35:GLN:HG3	9:CJ:36:VAL:H	1.81	0.44
10:CK:113:THR:HG23	10:CK:114:PRO:HD2	2.00	0.44
11:CL:85:ARG:HA	11:CL:93:ARG:HA	1.98	0.44
12:CM:10:ASP:C	12:CM:12:LYS:H	2.21	0.44
12:CM:63:VAL:HG13	12:CM:67:ASP:OD2	2.18	0.44
12:CM:79:LEU:C	12:CM:81:ASP:H	2.21	0.44
20:CO:19:ALA:O	20:CO:20:ASN:HB3	2.18	0.44
20:CO:25:THR:O	20:CO:26:GLU:C	2.56	0.44
14:CQ:80:LYS:HZ2	14:CQ:80:LYS:H	1.66	0.44
23:DB:1099:G:OP2	24:DI:2:LYS:O	2.35	0.44
23:DB:1173:U:H2'	23:DB:1174:U:C6	2.52	0.44
23:DB:1227:G:O2'	23:DB:1228:G:H5'	2.16	0.44
23:DB:1401:G:H2'	23:DB:1402:U:C6	2.53	0.44
23:DB:1528:A:H2'	23:DB:1529:G:O4'	2.18	0.44
23:DB:2376:A:H2'	23:DB:2377:A:O4'	2.17	0.44
23:DB:303:G:H2'	23:DB:304:U:H6	1.79	0.44
23:DB:460:A:H2'	23:DB:461:C:O4'	2.18	0.44
23:DB:805:G:N2	23:DB:829:A:OP1	2.51	0.44
23:DB:877:A:N6	23:DB:899:A:OP2	2.51	0.44
25:DC:116:GLN:O	25:DC:117:SER:HB3	2.17	0.44
29:DE:97:ASN:HD21	29:DE:100:MET:HG3	1.82	0.44
48:DG:91:VAL:C	48:DG:93:TYR:H	2.21	0.44
40:DH:72:ILE:CG2	40:DH:108:VAL:HG21	2.47	0.44
41:DJ:58:ASN:HA	41:DJ:127:GLY:HA2	2.00	0.44
41:DJ:64:VAL:CG2	41:DJ:68:LYS:HB2	2.48	0.44
37:DL:111:ILE:HA	37:DL:128:THR:HG23	2.00	0.44
38:DM:53:MET:SD	38:DM:63:ILE:HG21	2.57	0.44
43:DO:30:ARG:H	43:DO:97:PHE:HE2	1.64	0.44
43:DO:39:VAL:CG1	43:DO:48:LEU:HD12	2.42	0.44
28:DP:8:GLU:HA	28:DP:54:LEU:HD22	2.00	0.44
50:DT:54:GLU:O	50:DT:55:VAL:HB	2.18	0.44
50:DT:67:VAL:HG23	50:DT:75:GLY:O	2.17	0.44
30:DY:50:VAL:HB	30:DY:53:MET:HB3	2.00	0.44
1:AA:1041:G:O2'	1:AA:1042:A:H5'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1426:G:O2'	1:AA:1427:C:H5'	2.17	0.44
1:AA:319:G:O2'	1:AA:320:A:H5'	2.18	0.44
1:AA:113:G:O4'	1:AA:354:G:H4'	2.17	0.44
1:AA:412:A:H4'	1:AA:413:G:OP1	2.17	0.44
1:AA:456:A:H2'	1:AA:457:G:H5''	2.00	0.44
1:AA:472:U:H2'	1:AA:473:U:C6	2.52	0.44
1:AA:541:G:O2'	3:AD:39:GLN:HB3	2.17	0.44
1:AA:551:U:H2'	1:AA:552:U:C6	2.52	0.44
1:AA:679:C:O2'	1:AA:680:C:H5'	2.18	0.44
1:AA:696:A:H2'	1:AA:697:U:H6	1.83	0.44
1:AA:754:C:H3'	1:AA:754:C:O2	2.17	0.44
1:AA:763:G:H2'	1:AA:764:C:H6	1.82	0.44
1:AA:844:G:C2'	1:AA:845:A:H8	2.30	0.44
3:AD:154:VAL:O	3:AD:157:ALA:HB3	2.18	0.44
8:AI:46:VAL:HA	8:AI:49:GLN:CG	2.48	0.44
10:AK:113:THR:HG23	10:AK:114:PRO:HD2	2.00	0.44
10:AK:22:ILE:HD13	10:AK:95:THR:CG2	2.48	0.44
1:AA:552:U:H4'	11:AL:82:ARG:O	2.18	0.44
12:AM:100:ARG:C	12:AM:102:LYS:H	2.20	0.44
12:AM:33:LEU:HD12	12:AM:40:GLU:HG2	2.00	0.44
20:AO:77:ARG:HB3	20:AO:77:ARG:NH1	2.32	0.44
14:AQ:15:LYS:O	14:AQ:16:MET:C	2.56	0.44
14:AQ:3:LYS:N	14:AQ:3:LYS:HE2	2.32	0.44
15:AR:58:ILE:O	15:AR:61:ALA:HB3	2.18	0.44
16:AS:36:ARG:O	16:AS:69:LYS:HD2	2.17	0.44
31:B0:12:ARG:HD2	31:B0:16:ARG:NH1	2.33	0.44
22:BA:106:G:H2'	22:BA:107:G:C8	2.52	0.44
22:BA:73:A:H2'	22:BA:73:A:N3	2.33	0.44
22:BA:6:G:O2'	22:BA:7:G:H5'	2.18	0.44
23:BB:1030:C:O2'	23:BB:1031:G:H5'	2.17	0.44
23:BB:1059:G:H2'	23:BB:1060:U:C6	2.53	0.44
23:BB:1219:U:H2'	23:BB:1220:G:C8	2.53	0.44
23:BB:1369:G:O2'	23:BB:1370:C:H5'	2.17	0.44
23:BB:138:U:H1'	50:BT:1:MET:O	2.18	0.44
23:BB:1842:G:O2'	23:BB:1843:C:H5'	2.18	0.44
23:BB:2070:A:H2'	23:BB:2071:A:H8	1.80	0.44
23:BB:2153:C:H2'	23:BB:2154:A:C8	2.53	0.44
23:BB:2298:A:OP1	47:BF:70:ARG:HD3	2.18	0.44
23:BB:2648:G:H2'	23:BB:2649:C:C6	2.53	0.44
23:BB:265:A:O2'	23:BB:266:G:H4'	2.16	0.44
23:BB:2785:C:H2'	23:BB:2786:U:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:402:A:O2'	23:BB:403:U:H5'	2.17	0.44
23:BB:56:A:O2'	23:BB:57:C:H5'	2.18	0.44
23:BB:745:G:O2'	23:BB:748:G:H1'	2.17	0.44
23:BB:841:G:O2'	23:BB:842:U:H5'	2.18	0.44
29:BE:165:HIS:O	29:BE:167:VAL:N	2.47	0.44
47:BF:176:PHE:HB3	47:BF:177:ARG:H	1.67	0.44
48:BG:18:ILE:HA	48:BG:23:ILE:CG1	2.47	0.44
40:BH:29:PHE:O	40:BH:30:LEU:C	2.55	0.44
27:BK:18:ARG:O	27:BK:45:GLU:HB2	2.17	0.44
37:BL:121:THR:CB	37:BL:141:LYS:HB3	2.47	0.44
38:BM:110:GLU:H	38:BM:110:GLU:CD	2.21	0.44
28:BP:56:SER:O	28:BP:74:GLN:HA	2.17	0.44
44:BQ:57:ARG:C	44:BQ:59:LEU:N	2.71	0.44
44:BQ:65:ASN:O	44:BQ:69:ARG:HB2	2.18	0.44
39:BX:46:VAL:O	39:BX:50:VAL:HG23	2.18	0.44
51:BZ:32:ASN:HB3	51:BZ:34:HIS:NE2	2.33	0.44
1:CA:1005:A:O2'	1:CA:1037:C:H1'	2.18	0.44
1:CA:1091:U:H3'	6:CG:3:ARG:HH12	1.83	0.44
1:CA:131:A:H2'	1:CA:132:C:H6	1.82	0.44
1:CA:740:U:OP1	20:CO:38:HIS:HE1	2.00	0.44
1:CA:911:U:O2'	1:CA:912:C:H5'	2.17	0.44
18:CB:100:LEU:HD23	18:CB:174:GLU:C	2.38	0.44
18:CB:177:ASN:OD1	18:CB:178:LEU:HD23	2.18	0.44
18:CB:33:ALA:CB	18:CB:38:HIS:HA	2.48	0.44
18:CB:55:GLU:OE1	18:CB:58:LYS:HD2	2.17	0.44
2:CC:99:GLN:O	2:CC:100:ILE:HB	2.18	0.44
5:CF:3:HIS:CD2	5:CF:92:THR:HG23	2.53	0.44
5:CF:38:ARG:HH21	5:CF:63:ASN:ND2	2.16	0.44
6:CG:132:THR:O	6:CG:135:LYS:HB3	2.18	0.44
7:CH:43:GLY:HA2	7:CH:63:LYS:HZ1	1.82	0.44
8:CI:20:ILE:CD1	8:CI:62:LEU:HD12	2.47	0.44
1:CA:1367:C:H4'	9:CJ:50:THR:HG21	1.99	0.44
11:CL:42:LYS:HD2	11:CL:43:LYS:CG	2.48	0.44
21:CN:27:LYS:HG3	21:CN:28:ALA:N	2.26	0.44
14:CQ:30:HIS:CG	14:CQ:33:TYR:HB2	2.53	0.44
14:CQ:44:HIS:HB2	14:CQ:69:THR:O	2.18	0.44
15:CR:22:TYR:C	15:CR:24:ASP:H	2.22	0.44
17:CT:44:ALA:O	17:CT:48:LYS:HB3	2.18	0.44
36:D2:29:GLN:HE21	36:D2:33:ARG:HD2	1.83	0.44
23:DB:1145:C:O2'	23:DB:1146:C:H5'	2.16	0.44
23:DB:1457:U:H5'	23:DB:1458:U:OP1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1465:G:H2'	23:DB:1466:U:C6	2.53	0.44
23:DB:2276:G:O2'	23:DB:2277:G:H5'	2.18	0.44
23:DB:2771:C:H2'	23:DB:2772:C:C6	2.52	0.44
23:DB:538:A:H2'	23:DB:539:G:O4'	2.17	0.44
23:DB:704:G:H1'	23:DB:727:A:N6	2.31	0.44
23:DB:963:U:H2'	23:DB:964:C:C6	2.53	0.44
23:DB:969:G:O2'	23:DB:970:U:H5'	2.16	0.44
26:DD:32:ASN:HA	26:DD:51:THR:O	2.17	0.44
47:DF:33:ILE:CG2	47:DF:34:THR:N	2.81	0.44
48:DG:39:ALA:C	48:DG:54:ARG:HG3	2.38	0.44
24:DI:140:GLU:CD	24:DI:140:GLU:H	2.20	0.44
24:DI:70:THR:O	24:DI:70:THR:HG23	2.17	0.44
24:DI:57:VAL:HG23	24:DI:71:LYS:HZ1	1.81	0.44
24:DI:78:LEU:HD23	24:DI:81:LYS:HE2	2.00	0.44
41:DJ:45:THR:HG23	41:DJ:45:THR:O	2.17	0.44
41:DJ:14:ASP:O	41:DJ:53:TYR:HB2	2.17	0.44
38:DM:43:ALA:O	38:DM:46:ILE:HG12	2.17	0.44
43:DO:106:LEU:HG	43:DO:107:ALA:N	2.32	0.44
28:DP:3:ILE:C	28:DP:3:ILE:HD13	2.38	0.44
45:DS:108:SER:OG	45:DS:109:ASP:N	2.50	0.44
50:DT:39:THR:C	50:DT:41:ALA:H	2.21	0.44
35:DV:38:LEU:HG	35:DV:40:ILE:CD1	2.44	0.44
30:DY:21:ALA:O	30:DY:24:LEU:HB3	2.17	0.44
1:AA:216:U:H2'	1:AA:217:C:H6	1.83	0.44
1:AA:33:A:OP2	1:AA:398:U:H5'	2.18	0.44
1:AA:511:C:O2'	1:AA:512:U:H6	2.01	0.44
1:AA:737:C:H2'	1:AA:738:C:C6	2.53	0.44
1:AA:815:A:H4'	1:AA:817:C:C5	2.53	0.44
1:AA:13:U:C1'	1:AA:914:A:H5''	2.40	0.44
1:AA:996:A:N6	1:AA:1046:A:H1'	2.33	0.44
2:AC:47:ALA:C	2:AC:49:ALA:H	2.21	0.44
5:AF:18:VAL:HB	5:AF:19:PRO:CD	2.48	0.44
5:AF:92:THR:HG22	5:AF:93:LYS:N	2.31	0.44
7:AH:34:ALA:HB1	7:AH:109:VAL:HB	1.99	0.44
8:AI:85:ALA:O	8:AI:88:GLU:HB2	2.18	0.44
21:AN:55:SER:HB2	21:AN:58:ARG:HD2	2.00	0.44
22:BA:87:U:H2'	22:BA:88:C:H5''	1.99	0.44
23:BB:1292:G:H2'	23:BB:1293:C:C6	2.52	0.44
23:BB:1423:G:H2'	23:BB:1424:G:H8	1.83	0.44
23:BB:1486:U:H2'	23:BB:1487:U:C6	2.53	0.44
23:BB:1516:G:H2'	23:BB:1517:G:H8	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1635:A:H2'	23:BB:1636:U:O4'	2.17	0.44
23:BB:2273:A:H2'	23:BB:2274:A:C8	2.53	0.44
23:BB:2699:C:H2'	23:BB:2700:A:C8	2.53	0.44
23:BB:299:A:H2'	23:BB:300:A:C8	2.53	0.44
23:BB:305:C:O2'	23:BB:306:U:H5'	2.17	0.44
23:BB:485:C:HO2'	45:BS:60:HIS:CE1	2.35	0.44
23:BB:506:G:C5'	23:BB:509:C:H1'	2.46	0.44
23:BB:930:G:H1'	30:BY:24:LEU:HD11	2.00	0.44
23:BB:972:A:OP2	23:BB:974:G:H5''	2.18	0.44
23:BB:1797:G:O3'	25:BC:255:LYS:O	2.36	0.44
26:BD:34:VAL:O	26:BD:93:GLY:HA2	2.18	0.44
29:BE:115:GLN:C	29:BE:117:ARG:H	2.21	0.44
29:BE:6:LYS:HB3	29:BE:7:ASP:H	1.72	0.44
48:BG:96:ALA:CB	48:BG:103:ASN:HB3	2.41	0.44
41:BJ:83:GLY:O	41:BJ:84:ILE:C	2.57	0.44
27:BK:79:PHE:CD1	27:BK:79:PHE:N	2.83	0.44
37:BL:129:LYS:H	37:BL:129:LYS:HG2	1.56	0.44
43:BO:100:HIS:C	43:BO:104:GLN:HB2	2.38	0.44
28:BP:20:ARG:HB3	28:BP:23:ASP:OD2	2.17	0.44
28:BP:29:VAL:HA	28:BP:79:VAL:O	2.18	0.44
44:BQ:33:VAL:O	44:BQ:34:ALA:C	2.57	0.44
49:BR:34:GLU:HB3	49:BR:58:VAL:HG21	1.98	0.44
49:BR:19:THR:HA	49:BR:97:LYS:HA	2.00	0.44
45:BS:45:VAL:O	45:BS:48:LYS:HB3	2.18	0.44
50:BT:43:ILE:HG23	50:BT:58:VAL:HG21	1.99	0.44
23:BB:328:U:H4'	46:BU:65:GLN:CD	2.37	0.44
52:BW:32:ALA:O	52:BW:34:SER:N	2.49	0.44
1:CA:1348:U:O3'	8:CI:121:ARG:HG3	2.17	0.44
1:CA:153:C:H2'	1:CA:154:U:C6	2.53	0.44
1:CA:51:A:H4'	1:CA:52:C:OP2	2.18	0.44
1:CA:600:A:H2'	1:CA:601:G:C8	2.53	0.44
18:CB:169:HIS:CD2	18:CB:170:ILE:H	2.35	0.44
2:CC:19:SER:O	21:CN:93:PRO:HB3	2.18	0.44
8:CI:16:ALA:HA	8:CI:66:VAL:HA	1.98	0.44
11:CL:74:GLN:HB3	11:CL:75:GLU:H	1.48	0.44
21:CN:52:ARG:HA	21:CN:52:ARG:HD3	1.79	0.44
20:CO:7:ALA:O	20:CO:10:LYS:HB3	2.18	0.44
13:CP:67:ILE:HG12	13:CP:72:ALA:CB	2.48	0.44
14:CQ:8:GLN:HA	14:CQ:58:VAL:O	2.17	0.44
22:DA:85:G:H2'	22:DA:86:G:C8	2.52	0.44
23:DB:1044:C:O3'	23:DB:1047:G:H5'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:106:C:H2'	23:DB:107:G:C8	2.52	0.44
23:DB:1389:G:O2'	23:DB:1390:U:H5'	2.18	0.44
23:DB:1570:A:H2'	23:DB:1571:A:C8	2.52	0.44
23:DB:1684:G:H2'	23:DB:1685:C:C6	2.52	0.44
23:DB:1936:A:H2	23:DB:1943:U:O4	2.01	0.44
23:DB:1:G:H2'	23:DB:2:G:H8	1.79	0.44
23:DB:230:G:H2'	23:DB:231:A:C8	2.53	0.44
23:DB:2636:C:H2'	23:DB:2637:U:C6	2.52	0.44
23:DB:2666:C:O2	23:DB:2666:C:O4'	2.33	0.44
23:DB:745:G:O2'	23:DB:748:G:H1'	2.17	0.44
23:DB:861:A:H2'	23:DB:862:G:O4'	2.18	0.44
23:DB:948:C:H1'	23:DB:984:A:N3	2.33	0.44
25:DC:143:VAL:HA	25:DC:189:ALA:HB1	2.00	0.44
26:DD:122:VAL:HA	26:DD:127:PHE:H	1.83	0.44
29:DE:129:PRO:HB3	29:DE:159:LEU:HD23	1.99	0.44
29:DE:134:LEU:O	29:DE:138:LEU:HG	2.17	0.44
29:DE:165:HIS:O	29:DE:167:VAL:N	2.47	0.44
47:DF:11:VAL:HG12	47:DF:12:VAL:N	2.28	0.44
40:DH:5:LEU:HG	40:DH:17:ASP:O	2.18	0.44
27:DK:18:ARG:O	27:DK:45:GLU:HB2	2.18	0.44
27:DK:66:LYS:HD2	27:DK:81:GLY:N	2.32	0.44
27:DK:93:GLN:HB3	27:DK:94:PRO:HD2	1.99	0.44
42:DN:79:LEU:HD23	42:DN:83:LEU:HB3	2.00	0.44
22:DA:114:C:H1'	43:DO:47:VAL:HG21	2.00	0.44
43:DO:67:ASN:O	43:DO:69:ASP:N	2.51	0.44
49:DR:58:VAL:HG13	49:DR:59:ILE:N	2.33	0.44
45:DS:61:ASN:HB3	45:DS:62:ASP:H	1.62	0.44
46:DU:51:LEU:N	46:DU:53:GLN:NE2	2.66	0.44
1:AA:1238:A:H2	1:AA:1241:G:H21	1.63	0.43
1:AA:1390:U:O2'	1:AA:1391:U:H5'	2.18	0.43
1:AA:218:U:H2'	1:AA:219:U:O4'	2.17	0.43
1:AA:778:G:H2'	1:AA:779:C:H6	1.80	0.43
2:AC:156:LEU:HD11	2:AC:165:GLU:HB2	1.99	0.43
2:AC:57:GLU:CD	2:AC:64:ARG:HG2	2.38	0.43
3:AD:154:VAL:HG23	3:AD:155:LYS:N	2.33	0.43
8:AI:48:ARG:HA	8:AI:56:MET:SD	2.58	0.43
13:AP:68:SER:HB3	13:AP:71:VAL:HG12	1.99	0.43
16:AS:6:LYS:HD2	16:AS:6:LYS:N	2.32	0.43
17:AT:11:ILE:HG13	17:AT:11:ILE:H	1.66	0.43
31:B0:47:TYR:N	31:B0:47:TYR:CD1	2.86	0.43
32:B4:19:ARG:O	32:B4:21:GLY:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1076:C:H2'	23:BB:1077:A:H8	1.82	0.43
23:BB:1091:G:O2'	23:BB:1092:C:H5'	2.18	0.43
23:BB:1881:C:H2'	23:BB:1882:U:O4'	2.18	0.43
23:BB:2133:G:N3	23:BB:2133:G:C2'	2.74	0.43
23:BB:2336:A:O2'	23:BB:2337:G:P	2.76	0.43
23:BB:2533:U:H2'	23:BB:2534:A:O4'	2.17	0.43
23:BB:273:G:O2'	23:BB:274:C:H5'	2.18	0.43
23:BB:2821:A:H2'	23:BB:2822:G:C8	2.53	0.43
23:BB:297:G:OP1	46:BU:1:ALA:HB2	2.18	0.43
23:BB:714:U:H1'	23:BB:717:C:H5	1.82	0.43
25:BC:68:ARG:HE	25:BC:190:THR:HG22	1.83	0.43
26:BD:204:LYS:HB3	26:BD:204:LYS:NZ	2.33	0.43
26:BD:46:ARG:HG3	26:BD:46:ARG:NH1	2.33	0.43
23:BB:659:G:H21	29:BE:30:GLN:NE2	2.15	0.43
47:BF:148:VAL:O	47:BF:149:ARG:CB	2.65	0.43
47:BF:149:ARG:HA	47:BF:149:ARG:HD3	1.67	0.43
48:BG:91:VAL:C	48:BG:93:TYR:H	2.21	0.43
48:BG:93:TYR:O	48:BG:94:ARG:O	2.36	0.43
40:BH:66:ASN:O	40:BH:138:VAL:HG11	2.17	0.43
24:BI:102:ARG:HA	24:BI:105:LEU:HD12	2.00	0.43
24:BI:12:VAL:HG23	24:BI:41:PHE:CE2	2.53	0.43
41:BJ:3:THR:HB	41:BJ:44:TYR:CE1	2.52	0.43
27:BK:36:GLY:HA2	27:BK:62:VAL:O	2.18	0.43
37:BL:125:LEU:H	37:BL:143:GLU:HG3	1.83	0.43
37:BL:74:THR:HA	37:BL:107:PHE:O	2.17	0.43
37:BL:83:ALA:O	37:BL:85:VAL:N	2.49	0.43
38:BM:43:ALA:O	38:BM:46:ILE:HG12	2.18	0.43
42:BN:33:ILE:HD12	42:BN:34:ILE:N	2.33	0.43
43:BO:7:ARG:HD3	43:BO:97:PHE:CE1	2.53	0.43
44:BQ:104:ALA:C	44:BQ:106:THR:H	2.21	0.43
44:BQ:4:LYS:HE3	44:BQ:8:ILE:HD11	2.00	0.43
49:BR:59:ILE:HA	49:BR:101:ILE:H	1.83	0.43
45:BS:82:MET:HE2	45:BS:84:ARG:HH12	1.83	0.43
46:BU:23:LYS:HD2	46:BU:23:LYS:N	2.33	0.43
46:BU:66:VAL:C	46:BU:68:ASN:H	2.22	0.43
46:BU:86:PHE:CE1	46:BU:88:ASP:HB2	2.52	0.43
35:BV:29:ILE:HD12	35:BV:90:ASP:HA	2.00	0.43
23:BB:2336:A:H61	52:BW:40:ARG:HG3	1.82	0.43
1:CA:1314:C:H2'	1:CA:1315:U:C6	2.53	0.43
1:CA:1388:C:H2'	1:CA:1389:C:C6	2.53	0.43
1:CA:1392:G:C2'	1:CA:1393:U:H5'	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1509:C:O2'	1:CA:1510:C:H5'	2.18	0.43
1:CA:214:C:H2'	1:CA:215:C:C6	2.53	0.43
1:CA:184:G:H4'	1:CA:224:U:O3'	2.18	0.43
1:CA:509:A:N3	1:CA:543:U:O2'	2.51	0.43
1:CA:791:G:C6	1:CA:792:A:N7	2.86	0.43
1:CA:832:G:O2'	1:CA:833:G:H5'	2.18	0.43
1:CA:861:G:O2'	1:CA:862:C:H5'	2.18	0.43
18:CB:69:VAL:CG2	18:CB:162:VAL:HB	2.48	0.43
18:CB:162:VAL:HG13	18:CB:184:ALA:HB2	1.99	0.43
18:CB:69:VAL:HB	18:CB:162:VAL:HG23	1.99	0.43
2:CC:110:LEU:HD11	2:CC:143:LEU:O	2.18	0.43
2:CC:75:VAL:O	2:CC:75:VAL:HG12	2.17	0.43
2:CC:8:GLY:O	2:CC:11:LEU:HG	2.18	0.43
7:CH:74:ILE:HG12	7:CH:74:ILE:O	2.18	0.43
8:CI:32:ARG:CB	8:CI:36:GLN:HE21	2.29	0.43
12:CM:6:ILE:O	12:CM:7:ASN:C	2.57	0.43
17:CT:49:ALA:C	17:CT:52:GLU:HB3	2.38	0.43
22:DA:43:C:H5'	47:DF:62:GLN:CB	2.48	0.43
23:DB:1386:C:H2'	23:DB:1387:A:H8	1.83	0.43
23:DB:142:A:H2	50:DT:2:ILE:CG2	2.31	0.43
23:DB:2071:A:H2	23:DB:2440:C:N4	2.16	0.43
23:DB:2329:U:H2'	23:DB:2330:G:C8	2.53	0.43
23:DB:2425:A:H5''	23:DB:2426:A:H3'	1.99	0.43
23:DB:255:A:H2'	23:DB:256:A:O4'	2.18	0.43
23:DB:2604:U:O2'	23:DB:2605:U:H5'	2.18	0.43
23:DB:2734:A:C2'	23:DB:2735:G:H5'	2.45	0.43
23:DB:493:G:O2'	23:DB:494:G:H5'	2.17	0.43
23:DB:920:A:O2'	23:DB:921:C:H5'	2.18	0.43
25:DC:207:ALA:O	25:DC:208:GLY:C	2.56	0.43
25:DC:40:GLY:O	25:DC:53:ILE:HG23	2.17	0.43
26:DD:13:ARG:HH12	28:DP:74:GLN:NE2	2.16	0.43
29:DE:115:GLN:C	29:DE:117:ARG:H	2.20	0.43
29:DE:126:VAL:HG13	29:DE:156:ASN:HD22	1.83	0.43
29:DE:2:GLU:OE1	29:DE:13:THR:N	2.51	0.43
29:DE:58:LYS:HB2	29:DE:60:TRP:HD1	1.78	0.43
47:DF:69:ALA:O	47:DF:70:ARG:HB2	2.17	0.43
22:DA:43:C:C1'	47:DF:91:ARG:HD2	2.44	0.43
23:DB:2746:U:H4'	48:DG:137:LYS:HB2	1.98	0.43
40:DH:62:LEU:HG	40:DH:66:ASN:ND2	2.32	0.43
40:DH:90:LEU:HB2	40:DH:123:ARG:O	2.17	0.43
24:DI:21:PRO:HB2	24:DI:22:PRO:CD	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DK:112:PHE:O	27:DK:113:MET:C	2.56	0.43
37:DL:29:LYS:C	37:DL:31:GLY:H	2.21	0.43
38:DM:35:ALA:O	38:DM:36:VAL:HB	2.18	0.43
43:DO:9:ARG:O	43:DO:12:THR:HB	2.17	0.43
26:DD:15:PHE:H	28:DP:11:GLN:HE22	1.63	0.43
23:DB:518:G:H4'	45:DS:18:ARG:NH2	2.33	0.43
45:DS:24:ILE:O	45:DS:25:ARG:C	2.56	0.43
50:DT:4:GLU:C	50:DT:6:ARG:H	2.22	0.43
52:DW:31:LEU:O	52:DW:32:ALA:HB2	2.18	0.43
30:DY:4:ILE:HG12	30:DY:38:GLU:O	2.18	0.43
30:DY:4:ILE:HD12	30:DY:58:GLU:HG3	1.99	0.43
1:AA:1514:G:O2'	1:AA:1515:G:H5'	2.18	0.43
1:AA:254:G:H2'	1:AA:255:G:H8	1.83	0.43
1:AA:306:A:H2'	1:AA:307:C:H5'	1.99	0.43
1:AA:764:C:H2'	1:AA:765:G:C5'	2.47	0.43
1:AA:874:G:O2'	1:AA:875:U:H5'	2.17	0.43
1:AA:937:A:C2'	1:AA:938:A:H5'	2.49	0.43
1:AA:975:A:H4'	1:AA:976:G:OP2	2.18	0.43
18:AB:53:LEU:C	18:AB:55:GLU:N	2.72	0.43
6:AG:104:VAL:O	6:AG:108:ARG:HG3	2.18	0.43
1:AA:1368:A:OP2	8:AI:113:LYS:HD3	2.18	0.43
8:AI:10:ARG:HB2	8:AI:14:SER:O	2.17	0.43
8:AI:8:THR:O	8:AI:81:GLY:HA3	2.18	0.43
8:AI:95:SER:O	8:AI:99:LYS:HG3	2.17	0.43
14:AQ:37:ILE:HG22	14:AQ:38:LYS:N	2.34	0.43
19:AU:34:ARG:C	19:AU:34:ARG:NE	2.72	0.43
22:BA:25:U:H4'	22:BA:26:C:OP1	2.17	0.43
22:BA:43:C:H5'	47:BF:62:GLN:CB	2.48	0.43
23:BB:104:A:H2'	23:BB:105:C:H6	1.83	0.43
23:BB:1170:C:H2'	23:BB:1171:G:H8	1.81	0.43
23:BB:1334:G:O2'	23:BB:1335:C:H5'	2.18	0.43
23:BB:1949:G:H2'	23:BB:1950:G:C8	2.52	0.43
23:BB:2060:A:H2'	29:BE:63:LYS:CE	2.42	0.43
23:BB:2297:A:C2	23:BB:2320:U:H4'	2.52	0.43
23:BB:2606:C:O2'	23:BB:2607:G:H5'	2.18	0.43
23:BB:2659:G:H1'	23:BB:2662:A:N6	2.32	0.43
23:BB:2700:A:O2'	23:BB:2701:U:H5'	2.19	0.43
23:BB:287:G:O2'	23:BB:288:U:H5'	2.18	0.43
23:BB:346:A:H2'	23:BB:347:A:O4'	2.18	0.43
23:BB:776:G:H4'	23:BB:777:G:O5'	2.18	0.43
25:BC:161:VAL:HG12	25:BC:173:LEU:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:16:VAL:N	25:BC:203:VAL:HG11	2.28	0.43
25:BC:204:LEU:HD23	25:BC:209:ALA:CB	2.48	0.43
26:BD:187:LEU:HD12	26:BD:188:LEU:H	1.82	0.43
29:BE:52:VAL:CG1	29:BE:53:THR:H	2.22	0.43
47:BF:102:LEU:HD12	47:BF:103:ILE:HG13	2.00	0.43
47:BF:107:VAL:O	47:BF:110:ILE:HG22	2.17	0.43
47:BF:137:PHE:CD2	47:BF:137:PHE:N	2.87	0.43
47:BF:89:THR:OG1	47:BF:91:ARG:NH2	2.51	0.43
48:BG:10:VAL:HG23	48:BG:48:THR:HA	2.00	0.43
48:BG:40:VAL:HG13	48:BG:64:ALA:CB	2.48	0.43
38:BM:116:ALA:HA	38:BM:119:LEU:HD12	2.00	0.43
38:BM:31:PHE:CZ	38:BM:110:GLU:HB3	2.52	0.43
38:BM:43:ALA:HB2	38:BM:92:TRP:O	2.18	0.43
42:BN:37:THR:CG2	42:BN:39:PRO:HD2	2.35	0.43
44:BQ:10:ARG:NH1	44:BQ:10:ARG:HB2	2.33	0.43
44:BQ:7:VAL:CG2	44:BQ:8:ILE:N	2.81	0.43
45:BS:41:LYS:O	45:BS:43:ALA:N	2.51	0.43
46:BU:40:LEU:HA	46:BU:61:GLU:HA	1.99	0.43
35:BV:56:PHE:O	35:BV:61:LEU:HD21	2.18	0.43
1:CA:991:U:C4	1:CA:1212:U:H4'	2.52	0.43
1:CA:1309:G:H2'	1:CA:1310:G:H8	1.84	0.43
1:CA:586:C:H5''	7:CH:81:GLY:HA2	2.00	0.43
1:CA:845:A:H5''	1:CA:846:G:H8	1.80	0.43
1:CA:852:G:H2'	1:CA:853:C:H6	1.82	0.43
18:CB:212:TYR:HA	18:CB:215:ALA:CB	2.44	0.43
18:CB:216:VAL:CG2	18:CB:217:ALA:N	2.79	0.43
2:CC:54:ILE:O	2:CC:54:ILE:HG23	2.18	0.43
3:CD:138:PRO:C	3:CD:140:ASP:H	2.21	0.43
5:CF:36:ILE:HG12	5:CF:64:VAL:HG22	2.00	0.43
6:CG:43:TYR:O	6:CG:46:LEU:HB3	2.18	0.43
12:CM:95:PRO:HA	12:CM:108:ARG:HG2	1.99	0.43
12:CM:89:ARG:HD2	12:CM:95:PRO:O	2.18	0.43
1:CA:618:C:H1'	13:CP:14:ARG:NH1	2.32	0.43
15:CR:42:ARG:C	15:CR:44:THR:H	2.20	0.43
31:D0:2:VAL:HG12	31:D0:3:GLN:H	1.83	0.43
33:D1:18:HIS:CE1	33:D1:40:PRO:HD2	2.53	0.43
36:D2:20:ALA:C	36:D2:22:MET:N	2.71	0.43
23:DB:1281:G:H2'	23:DB:1282:U:O4'	2.19	0.43
23:DB:1338:G:C2'	23:DB:1339:G:H5'	2.48	0.43
23:DB:1348:C:H5'	23:DB:1349:C:OP2	2.18	0.43
23:DB:151:C:H2'	23:DB:152:A:H8	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2271:G:O2'	23:DB:2272:U:H5'	2.18	0.43
23:DB:2305:U:H2'	23:DB:2306:C:O4'	2.17	0.43
23:DB:2770:G:O5'	23:DB:2770:G:H8	2.02	0.43
23:DB:459:U:O2'	23:DB:460:A:H5'	2.18	0.43
23:DB:754:U:H2'	23:DB:755:U:H6	1.83	0.43
23:DB:798:G:O2'	23:DB:799:G:H5'	2.18	0.43
25:DC:61:TYR:HA	25:DC:85:ASN:ND2	2.33	0.43
26:DD:168:GLU:O	26:DD:169:ARG:C	2.57	0.43
47:DF:126:ASN:HD22	47:DF:156:THR:HA	1.83	0.43
47:DF:146:ASP:O	47:DF:147:ARG:C	2.56	0.43
47:DF:34:THR:OG1	47:DF:154:THR:HB	2.18	0.43
40:DH:28:ASN:HD21	51:DZ:36:HIS:CE1	2.35	0.43
37:DL:19:LEU:CD2	37:DL:27:LEU:HD23	2.48	0.43
38:DM:100:LYS:HD2	38:DM:101:VAL:H	1.83	0.43
38:DM:59:ARG:HH11	38:DM:60:GLN:HB3	1.82	0.43
27:DK:76:VAL:O	28:DP:71:ARG:HA	2.18	0.43
44:DQ:63:ARG:HH22	44:DQ:96:ASP:N	2.16	0.43
45:DS:29:VAL:HG23	45:DS:70:LYS:HA	1.99	0.43
35:DV:12:GLN:O	35:DV:16:ALA:HB3	2.18	0.43
51:DZ:53:ALA:O	51:DZ:55:GLY:N	2.45	0.43
51:DZ:53:ALA:O	51:DZ:54:LYS:HB3	2.17	0.43
51:DZ:68:LEU:O	51:DZ:69:ALA:C	2.57	0.43
1:AA:1126:U:HO2'	1:AA:1280:A:H2'	1.83	0.43
1:AA:1471:U:O2'	1:AA:1472:U:H5'	2.18	0.43
1:AA:1473:G:H2'	1:AA:1474:U:C6	2.53	0.43
1:AA:5:U:H1'	1:AA:6:G:N1	2.33	0.43
18:AB:44:LYS:C	18:AB:47:PRO:HD2	2.38	0.43
2:AC:181:ILE:N	2:AC:181:ILE:CD1	2.79	0.43
3:AD:58:GLN:HA	3:AD:58:GLN:OE1	2.18	0.43
8:AI:126:PHE:H	8:AI:126:PHE:HD2	1.67	0.43
8:AI:63:TYR:C	8:AI:64:ILE:HD12	2.39	0.43
8:AI:16:ALA:HA	8:AI:66:VAL:HA	2.00	0.43
20:AO:43:PHE:HE2	20:AO:53:ARG:HD3	1.83	0.43
13:AP:71:VAL:HG13	13:AP:72:ALA:H	1.82	0.43
23:BB:2285:C:OP2	33:B1:5:ARG:HD3	2.19	0.43
34:B3:16:THR:C	34:B3:18:LYS:H	2.21	0.43
23:BB:1047:G:H1'	23:BB:1111:A:N6	2.32	0.43
23:BB:1055:G:HO2'	23:BB:1085:A:H2	1.65	0.43
23:BB:1190:G:OP1	37:BL:32:GLY:HA2	2.18	0.43
23:BB:1459:G:H3'	23:BB:1460:U:H4'	2.00	0.43
23:BB:1682:G:H2'	23:BB:1683:U:C6	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2376:A:H2'	23:BB:2377:A:O4'	2.18	0.43
23:BB:2693:G:H2'	23:BB:2694:G:C8	2.51	0.43
23:BB:2834:G:H1'	23:BB:2883:A:H61	1.81	0.43
23:BB:483:A:H2'	23:BB:484:C:O4'	2.18	0.43
23:BB:574:A:H4'	23:BB:575:A:H5''	1.99	0.43
23:BB:685:A:H1'	23:BB:688:U:O4	2.18	0.43
23:BB:859:G:O2'	23:BB:860:U:OP2	2.35	0.43
23:BB:936:A:H2'	23:BB:937:C:H6	1.81	0.43
26:BD:109:VAL:HG11	26:BD:193:VAL:HG11	2.00	0.43
26:BD:38:LYS:CB	26:BD:47:ALA:HB3	2.49	0.43
29:BE:114:ARG:HG3	29:BE:114:ARG:HH11	1.83	0.43
29:BE:62:GLN:CG	29:BE:63:LYS:N	2.81	0.43
47:BF:151:LEU:CD1	47:BF:153:ILE:HG13	2.48	0.43
48:BG:10:VAL:HB	48:BG:47:ASN:O	2.18	0.43
41:BJ:25:LEU:HB2	41:BJ:62:VAL:HG22	2.00	0.43
23:BB:2276:G:O3'	38:BM:86:LYS:HB2	2.19	0.43
42:BN:29:VAL:O	42:BN:78:LYS:HG2	2.17	0.43
49:BR:63:VAL:HA	49:BR:95:ASP:O	2.18	0.43
50:BT:4:GLU:C	50:BT:6:ARG:H	2.20	0.43
39:BX:19:LEU:O	39:BX:24:GLU:HB2	2.18	0.43
30:BY:50:VAL:HB	30:BY:53:MET:HB3	1.99	0.43
1:CA:1248:A:H2'	1:CA:1249:C:H6	1.83	0.43
1:CA:1342:C:H1'	8:CI:125:GLN:OE1	2.19	0.43
1:CA:1476:A:H2'	1:CA:1477:U:H6	1.83	0.43
2:CC:129:PHE:CG	2:CC:130:ARG:N	2.86	0.43
5:CF:10:VAL:HA	5:CF:84:VAL:HA	2.00	0.43
8:CI:25:GLY:HA3	8:CI:57:VAL:HA	2.00	0.43
9:CJ:71:LEU:N	9:CJ:71:LEU:HD23	2.32	0.43
10:CK:111:ASP:CB	19:CU:19:LYS:HE2	2.48	0.43
12:CM:89:ARG:NH1	12:CM:101:THR:HG21	2.33	0.43
9:CJ:66:GLU:HG2	21:CN:98:ALA:HA	2.00	0.43
20:CO:5:THR:O	20:CO:8:THR:HB	2.18	0.43
17:CT:38:ILE:CD1	17:CT:82:ILE:HA	2.48	0.43
17:CT:59:ARG:HB2	17:CT:59:ARG:HH11	1.83	0.43
31:D0:41:HIS:O	31:D0:42:ILE:O	2.37	0.43
22:DA:67:G:O2'	22:DA:68:C:H5'	2.17	0.43
23:DB:1692:U:H2'	23:DB:1694:C:C4	2.54	0.43
23:DB:1712:U:H3'	23:DB:1713:A:H2'	2.00	0.43
23:DB:2026:U:H2'	23:DB:2027:G:H8	1.83	0.43
23:DB:213:A:O2'	23:DB:214:G:H5'	2.18	0.43
23:DB:2527:C:H2'	23:DB:2528:U:C6	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2630:G:H2'	23:DB:2631:G:C8	2.52	0.43
23:DB:2893:A:H4'	23:DB:2894:G:C5'	2.49	0.43
23:DB:899:A:H5'	23:DB:900:A:OP2	2.18	0.43
23:DB:970:U:O2'	23:DB:984:A:H4'	2.18	0.43
26:DD:121:THR:C	26:DD:123:LYS:N	2.71	0.43
47:DF:129:MET:N	47:DF:129:MET:SD	2.92	0.43
23:DB:2529:G:H4'	48:DG:174:LYS:HG3	1.98	0.43
48:DG:10:VAL:HG23	48:DG:48:THR:HA	2.00	0.43
48:DG:93:TYR:CD2	48:DG:93:TYR:N	2.87	0.43
40:DH:104:THR:HA	40:DH:108:VAL:O	2.19	0.43
40:DH:5:LEU:CD1	40:DH:17:ASP:HB3	2.39	0.43
27:DK:20:MET:C	27:DK:41:ILE:HG13	2.39	0.43
37:DL:85:VAL:O	37:DL:85:VAL:HG22	2.18	0.43
43:DO:115:LEU:H	43:DO:115:LEU:HD22	1.83	0.43
28:DP:50:ARG:HB3	28:DP:57:ALA:O	2.18	0.43
28:DP:77:SER:O	28:DP:80:VAL:HG12	2.18	0.43
44:DQ:7:VAL:CG2	44:DQ:8:ILE:N	2.81	0.43
45:DS:35:ILE:HG22	45:DS:39:THR:CG2	2.46	0.43
45:DS:82:MET:HE2	45:DS:84:ARG:HH12	1.83	0.43
50:DT:23:ALA:C	50:DT:25:GLU:H	2.20	0.43
46:DU:16:LYS:HB3	46:DU:17:ASP:H	1.68	0.43
1:AA:1146:A:N3	1:AA:1146:A:H2'	2.33	0.43
1:AA:1152:A:H2'	1:AA:1153:G:C8	2.53	0.43
1:AA:1530:G:H2'	1:AA:1531:A:C8	2.54	0.43
1:AA:777:A:H2'	1:AA:778:G:C8	2.53	0.43
1:AA:840:C:H2'	1:AA:843:U:OP2	2.19	0.43
18:AB:22:TRP:HA	18:AB:189:ASN:N	2.33	0.43
2:AC:55:VAL:O	2:AC:65:VAL:HA	2.18	0.43
4:AE:93:VAL:HG12	4:AE:94:PHE:N	2.33	0.43
5:AF:16:GLU:CD	5:AF:16:GLU:H	2.20	0.43
7:AH:17:GLN:HE21	7:AH:62:LEU:HG	1.84	0.43
7:AH:63:LYS:CG	7:AH:70:VAL:HG21	2.47	0.43
10:AK:17:ASP:C	10:AK:36:ARG:HH12	2.20	0.43
11:AL:106:VAL:HA	11:AL:107:LYS:HZ3	1.82	0.43
12:AM:56:ARG:HG3	12:AM:56:ARG:NH1	2.31	0.43
21:AN:58:ARG:N	21:AN:59:GLN:HE21	2.14	0.43
14:AQ:40:THR:HG22	14:AQ:41:THR:N	2.32	0.43
23:BB:2526:G:H1'	32:B4:1:MET:H1	1.81	0.43
23:BB:1058:U:O2'	23:BB:1059:G:H5'	2.19	0.43
23:BB:118:A:H5'	23:BB:119:A:C8	2.46	0.43
23:BB:1296:G:O2'	23:BB:1297:C:H5'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1348:C:H5'	23:BB:1349:C:OP2	2.18	0.43
23:BB:1353:A:O2'	23:BB:1354:A:H5'	2.18	0.43
23:BB:1360:G:C2'	23:BB:1361:G:H5'	2.47	0.43
23:BB:1541:C:O2'	23:BB:1542:U:H5'	2.19	0.43
23:BB:1839:G:C8	23:BB:1927:A:H1'	2.53	0.43
23:BB:1920:C:H2'	23:BB:1921:G:H8	1.82	0.43
23:BB:2271:G:O2'	23:BB:2272:U:H5'	2.18	0.43
23:BB:2800:A:O5'	23:BB:2800:A:H8	2.01	0.43
23:BB:546:U:H5''	23:BB:546:U:O2	2.18	0.43
23:BB:915:C:H3'	23:BB:916:G:C8	2.52	0.43
25:BC:33:LEU:O	25:BC:34:GLU:HB3	2.18	0.43
26:BD:187:LEU:O	26:BD:188:LEU:HD23	2.18	0.43
26:BD:36:GLN:O	26:BD:36:GLN:HG3	2.17	0.43
26:BD:62:LYS:CB	26:BD:63:PRO:HD3	2.47	0.43
29:BE:60:TRP:CZ3	29:BE:69:ARG:HA	2.53	0.43
47:BF:110:ILE:HG22	47:BF:113:PHE:HB3	2.00	0.43
47:BF:99:PHE:O	47:BF:102:LEU:HB3	2.18	0.43
48:BG:77:GLY:HA3	48:BG:135:ALA:O	2.18	0.43
48:BG:6:ALA:HA	48:BG:7:PRO:HD3	1.82	0.43
24:BI:138:VAL:HG12	24:BI:139:VAL:N	2.33	0.43
41:BJ:73:VAL:HG23	41:BJ:74:TYR:H	1.83	0.43
38:BM:41:LEU:HB3	38:BM:46:ILE:HG21	1.99	0.43
43:BO:66:GLY:C	43:BO:68:LYS:H	2.21	0.43
49:BR:26:ASP:O	49:BR:27:ILE:HG12	2.18	0.43
45:BS:24:ILE:O	45:BS:25:ARG:C	2.56	0.43
45:BS:33:LEU:O	45:BS:37:THR:HG23	2.17	0.43
50:BT:69:ARG:NE	50:BT:69:ARG:HA	2.26	0.43
46:BU:41:VAL:HG22	46:BU:60:LYS:O	2.17	0.43
35:BV:38:LEU:CG	35:BV:40:ILE:HG23	2.47	0.43
1:CA:1096:C:O2'	1:CA:1097:C:H5'	2.18	0.43
1:CA:1179:A:H2'	1:CA:1180:A:O4'	2.18	0.43
1:CA:1312:G:H2'	1:CA:1313:U:H6	1.83	0.43
1:CA:216:U:H2'	1:CA:217:C:H6	1.82	0.43
1:CA:81:A:O2'	1:CA:82:G:H5'	2.18	0.43
1:CA:98:A:O2'	1:CA:99:C:H5'	2.19	0.43
2:CC:106:ARG:HD2	2:CC:106:ARG:C	2.38	0.43
2:CC:155:ARG:NE	2:CC:159:ALA:O	2.51	0.43
3:CD:96:ARG:HH12	3:CD:133:SER:HA	1.83	0.43
5:CF:15:SER:HA	5:CF:18:VAL:HG23	1.99	0.43
1:CA:878:A:C5'	7:CH:80:PRO:HG2	2.48	0.43
8:CI:7:GLY:HA3	8:CI:81:GLY:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CK:83:VAL:HB	10:CK:109:ILE:HG12	2.00	0.43
10:CK:22:ILE:HD13	10:CK:95:THR:CG2	2.48	0.43
11:CL:24:GLU:C	11:CL:26:CYS:N	2.70	0.43
12:CM:33:LEU:HB3	12:CM:38:ILE:O	2.18	0.43
12:CM:10:ASP:CA	12:CM:44:ILE:HD13	2.48	0.43
17:CT:8:LYS:HE3	17:CT:12:GLN:OE1	2.19	0.43
31:D0:27:LEU:HD22	31:D0:37:HIS:C	2.39	0.43
23:DB:2418:A:OP1	34:D3:44:ARG:NH1	2.52	0.43
22:DA:112:G:H2'	22:DA:113:C:H6	1.83	0.43
23:DB:1142:A:H4'	41:DJ:27:ARG:NH2	2.31	0.43
23:DB:1183:U:H2'	23:DB:1184:U:H6	1.83	0.43
23:DB:1316:U:O2'	23:DB:1317:G:H5'	2.18	0.43
23:DB:1473:G:O2'	23:DB:1474:U:H5'	2.18	0.43
23:DB:1516:G:H2'	23:DB:1517:G:H8	1.84	0.43
23:DB:1592:C:H2'	23:DB:1593:A:C8	2.54	0.43
23:DB:192:C:H2'	23:DB:193:U:H5'	2.00	0.43
23:DB:662:G:O2'	23:DB:663:G:H5'	2.18	0.43
23:DB:870:U:C2'	23:DB:871:U:H5''	2.48	0.43
25:DC:33:LEU:O	25:DC:34:GLU:HB3	2.18	0.43
25:DC:54:GLY:O	25:DC:214:GLY:HA2	2.18	0.43
48:DG:134:GLY:HA3	48:DG:140:ILE:HG21	1.99	0.43
37:DL:23:ILE:H	37:DL:23:ILE:CD1	2.10	0.43
38:DM:41:LEU:O	38:DM:94:ALA:N	2.49	0.43
43:DO:100:HIS:HA	43:DO:104:GLN:CG	2.48	0.43
45:DS:57:ASN:ND2	45:DS:57:ASN:N	2.67	0.43
50:DT:8:LEU:O	50:DT:9:LYS:HE2	2.19	0.43
46:DU:13:LEU:HA	46:DU:18:LYS:NZ	2.34	0.43
46:DU:3:LYS:HD3	46:DU:82:VAL:CG2	2.47	0.43
30:DY:47:ILE:HG21	30:DY:56:VAL:HG21	1.98	0.43
1:AA:731:G:O2'	1:AA:732:C:H5'	2.18	0.43
1:AA:792:A:N3	1:AA:794:A:C5	2.86	0.43
1:AA:862:C:O2'	1:AA:863:U:H5'	2.18	0.43
2:AC:13:ILE:HD13	2:AC:13:ILE:N	2.34	0.43
3:AD:191:SER:O	3:AD:192:ALA:HB2	2.19	0.43
3:AD:49:ASP:OD2	4:AE:111:ARG:HD2	2.18	0.43
6:AG:31:VAL:HG13	6:AG:32:ASP:OD1	2.19	0.43
21:AN:20:PHE:HE1	21:AN:51:PRO:HG3	1.84	0.43
19:AU:13:VAL:CG2	19:AU:15:LEU:HD23	2.48	0.43
22:BA:69:G:H2'	22:BA:70:C:H5'	2.00	0.43
23:BB:100:U:C6	23:BB:100:U:O5'	2.71	0.43
23:BB:1053:C:H6	23:BB:1053:C:O5'	2.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1281:G:H2'	23:BB:1282:U:O4'	2.18	0.43
23:BB:1359:A:H2'	23:BB:1360:G:O4'	2.17	0.43
23:BB:1465:G:H2'	23:BB:1466:U:C6	2.53	0.43
23:BB:1671:U:H2'	23:BB:1673:G:OP2	2.18	0.43
23:BB:1726:C:H2'	23:BB:1727:C:C6	2.53	0.43
23:BB:1747:U:H2'	23:BB:1748:C:C6	2.54	0.43
23:BB:1873:G:H2'	23:BB:1874:C:H6	1.83	0.43
23:BB:2626:C:H2'	23:BB:2627:G:H8	1.81	0.43
23:BB:2771:C:H2'	23:BB:2772:C:C6	2.53	0.43
23:BB:2897:U:H2'	23:BB:2898:U:H6	1.84	0.43
23:BB:443:A:N3	23:BB:443:A:H2'	2.33	0.43
23:BB:538:A:H2'	23:BB:539:G:O4'	2.18	0.43
23:BB:664:G:O2'	23:BB:665:U:H5'	2.18	0.43
23:BB:803:U:O2'	23:BB:804:A:H5'	2.18	0.43
23:BB:917:A:H2'	23:BB:918:A:O4'	2.18	0.43
25:BC:163:ILE:HG22	25:BC:164:VAL:H	1.83	0.43
25:BC:185:ALA:C	25:BC:187:CYS:H	2.22	0.43
26:BD:202:ILE:HG22	26:BD:204:LYS:HZ2	1.83	0.43
47:BF:32:LYS:N	47:BF:95:MET:SD	2.92	0.43
48:BG:10:VAL:O	48:BG:10:VAL:HG12	2.19	0.43
40:BH:108:VAL:HG12	40:BH:109:GLU:H	1.83	0.43
40:BH:112:LYS:HE3	40:BH:112:LYS:C	2.39	0.43
40:BH:26:ALA:O	40:BH:28:ASN:N	2.51	0.43
24:BI:85:ILE:CD1	24:BI:137:LEU:HD21	2.47	0.43
24:BI:79:LEU:HD23	24:BI:108:ILE:CD1	2.49	0.43
38:BM:82:MET:HE3	38:BM:83:GLY:H	1.84	0.43
42:BN:71:ARG:HG2	42:BN:71:ARG:HH21	1.83	0.43
28:BP:6:GLN:HE21	28:BP:7:LEU:HD12	1.84	0.43
44:BQ:82:LEU:HA	44:BQ:85:ALA:HB3	2.01	0.43
49:BR:98:ILE:HG22	49:BR:99:THR:N	2.33	0.43
46:BU:34:ILE:HG23	46:BU:63:ALA:HB2	2.01	0.43
23:BB:2264:C:H41	52:BW:11:ASN:HD22	1.67	0.43
30:BY:12:ALA:HB2	30:BY:53:MET:SD	2.58	0.43
30:BY:47:ILE:HG21	30:BY:56:VAL:HG21	2.00	0.43
1:CA:1379:G:N7	6:CG:2:ARG:NH1	2.66	0.43
1:CA:33:A:OP2	1:CA:398:U:H5'	2.17	0.43
1:CA:202:G:N2	1:CA:465:A:H61	2.12	0.43
1:CA:490:C:H2'	1:CA:491:G:O4'	2.18	0.43
1:CA:629:A:H2'	1:CA:630:A:O4'	2.18	0.43
1:CA:702:A:N6	23:DB:1848:A:N1	2.66	0.43
1:CA:865:A:H5'	1:CA:1078:U:C4	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CB:204:ASP:O	18:CB:205:ALA:HB3	2.18	0.43
1:CA:1206:G:C4'	2:CC:192:TYR:HA	2.41	0.43
3:CD:151:GLN:O	3:CD:154:VAL:HG22	2.18	0.43
4:CE:59:ILE:H	4:CE:59:ILE:HG13	1.56	0.43
5:CF:9:MET:HG2	5:CF:9:MET:H	1.73	0.43
7:CH:50:VAL:HG13	7:CH:50:VAL:O	2.18	0.43
7:CH:63:LYS:HG2	7:CH:70:VAL:CG2	2.47	0.43
1:CA:552:U:H4'	11:CL:82:ARG:O	2.17	0.43
12:CM:87:GLY:O	12:CM:91:ARG:HG3	2.18	0.43
20:CO:36:ILE:HD12	20:CO:60:VAL:CG2	2.46	0.43
19:CU:47:ALA:O	19:CU:51:ALA:N	2.51	0.43
31:D0:18:HIS:ND1	31:D0:18:HIS:N	2.67	0.43
23:DB:2756:U:OP2	32:D4:19:ARG:HG2	2.19	0.43
23:DB:1080:A:O2'	23:DB:1081:U:H5'	2.18	0.43
23:DB:1187:G:H5''	49:DR:83:TYR:CE2	2.52	0.43
23:DB:1330:C:H2'	23:DB:1331:G:C8	2.53	0.43
23:DB:1518:C:H2'	23:DB:1519:G:H8	1.84	0.43
23:DB:1568:G:O3'	25:DC:58:LYS:HD3	2.18	0.43
23:DB:1589:U:H2'	23:DB:1590:A:C8	2.54	0.43
23:DB:1625:C:H3'	23:DB:1626:A:C8	2.54	0.43
23:DB:2466:C:OP1	32:D4:4:ARG:HB3	2.19	0.43
23:DB:195:A:H1'	23:DB:250:G:H21	1.84	0.43
23:DB:2660:A:H2'	23:DB:2661:G:O4'	2.19	0.43
23:DB:2900:A:H2'	23:DB:2901:C:C6	2.53	0.43
23:DB:308:G:O4'	23:DB:501:A:H5'	2.18	0.43
23:DB:307:G:N2	23:DB:309:A:H3'	2.34	0.43
23:DB:528:A:H2'	23:DB:529:A:H5''	2.00	0.43
25:DC:163:ILE:HG22	25:DC:164:VAL:H	1.84	0.43
40:DH:69:ALA:HA	40:DH:140:ALA:CB	2.48	0.43
24:DI:59:THR:O	24:DI:59:THR:HG23	2.19	0.43
41:DJ:19:ASP:OD2	41:DJ:58:ASN:N	2.52	0.43
27:DK:62:VAL:HG12	27:DK:63:VAL:H	1.84	0.43
37:DL:80:SER:HB3	37:DL:115:GLU:CD	2.39	0.43
23:DB:826:U:O2'	37:DL:53:GLY:HA3	2.19	0.43
49:DR:55:ASP:CG	49:DR:56:GLY:H	2.21	0.43
45:DS:29:VAL:O	45:DS:32:ALA:HB3	2.18	0.43
45:DS:32:ALA:O	45:DS:35:ILE:HG12	2.19	0.43
35:DV:75:GLN:HA	35:DV:75:GLN:OE1	2.17	0.43
35:DV:29:ILE:HD12	35:DV:90:ASP:HA	2.00	0.43
39:DX:25:GLN:HB2	39:DX:46:VAL:HG11	2.00	0.43
1:AA:1057:G:O2'	1:AA:1058:G:H5'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:112:G:O2'	1:AA:113:G:H5'	2.18	0.43
1:AA:1054:C:H1'	1:AA:1196:A:C5	2.54	0.43
1:AA:1358:U:H5''	21:AN:72:PHE:O	2.19	0.43
1:AA:286:C:H2'	1:AA:287:U:H6	1.83	0.43
1:AA:337:G:O2'	1:AA:338:A:H5'	2.19	0.43
1:AA:430:A:OP2	3:AD:6:PRO:HA	2.17	0.43
1:AA:537:G:H2'	1:AA:538:G:C8	2.53	0.43
1:AA:634:C:O5'	1:AA:634:C:H6	2.01	0.43
1:AA:709:U:H2'	1:AA:710:G:H8	1.83	0.43
1:AA:884:U:H4'	1:AA:885:G:H5''	2.01	0.43
1:AA:937:A:H2'	1:AA:938:A:H5'	2.00	0.43
1:AA:959:A:H2	1:AA:1221:G:N3	2.17	0.43
18:AB:163:ILE:CG2	18:AB:164:ASP:N	2.81	0.43
6:AG:71:THR:HG22	6:AG:141:HIS:CE1	2.53	0.43
6:AG:73:GLU:C	6:AG:87:PRO:HA	2.38	0.43
7:AH:50:VAL:O	7:AH:50:VAL:HG13	2.19	0.43
8:AI:71:ILE:CD1	8:AI:71:ILE:H	2.28	0.43
13:AP:3:THR:HG22	13:AP:66:THR:HB	2.00	0.43
16:AS:44:ILE:HA	16:AS:61:VAL:CB	2.48	0.43
36:B2:1:MET:HG2	36:B2:2:LYS:H	1.83	0.43
22:BA:102:G:H2'	22:BA:103:U:C6	2.54	0.43
23:BB:1526:C:O2'	23:BB:1527:G:H5'	2.17	0.43
23:BB:1271:G:OP2	23:BB:1648:U:OP1	2.37	0.43
23:BB:1695:G:H1'	25:BC:7:PRO:O	2.17	0.43
23:BB:1824:G:OP2	25:BC:52:HIS:CE1	2.72	0.43
1:AA:1492:A:C4	23:BB:1913:A:N1	2.87	0.43
23:BB:1914:C:O2	23:BB:1914:C:C3'	2.67	0.43
1:AA:1483:A:H1'	23:BB:1948:G:H1'	2.00	0.43
23:BB:2361:G:O2'	23:BB:2362:C:H5'	2.18	0.43
23:BB:2599:G:H2'	23:BB:2600:A:C8	2.54	0.43
23:BB:2774:C:OP1	26:BD:169:ARG:HG3	2.18	0.43
23:BB:2868:A:H2'	23:BB:2869:G:C8	2.53	0.43
23:BB:372:G:O2'	23:BB:373:U:P	2.77	0.43
23:BB:451:U:C2	23:BB:453:A:N7	2.86	0.43
23:BB:560:C:H3'	23:BB:561:G:H8	1.82	0.43
23:BB:635:C:H2'	23:BB:636:G:C8	2.54	0.43
23:BB:870:U:C2'	23:BB:871:U:H5''	2.48	0.43
23:BB:969:G:O2'	23:BB:970:U:H5'	2.19	0.43
25:BC:140:VAL:CG1	25:BC:141:HIS:H	2.22	0.43
25:BC:23:LEU:HA	25:BC:23:LEU:HD12	1.86	0.43
48:BG:64:ALA:O	48:BG:67:ALA:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BG:84:LYS:CG	48:BG:85:LYS:H	2.19	0.43
48:BG:88:LEU:HD13	48:BG:93:TYR:HB2	2.00	0.43
40:BH:89:LYS:HB3	40:BH:90:LEU:H	1.48	0.43
24:BI:56:VAL:HG13	24:BI:58:ILE:HD11	2.01	0.43
41:BJ:131:ASN:O	41:BJ:132:HIS:ND1	2.52	0.43
27:BK:59:LYS:HD2	27:BK:89:ASN:ND2	2.33	0.43
37:BL:78:ARG:CB	37:BL:113:ALA:HB2	2.39	0.43
37:BL:120:VAL:HG12	37:BL:121:THR:N	2.34	0.43
28:BP:62:LYS:HB3	28:BP:69:VAL:HG22	1.99	0.43
35:BV:29:ILE:HG13	35:BV:88:HIS:HE1	1.83	0.43
51:BZ:40:VAL:O	51:BZ:42:SER:N	2.47	0.43
1:CA:1448:C:O2'	1:CA:1449:C:H5'	2.18	0.43
1:CA:254:G:H2'	1:CA:255:G:H8	1.83	0.43
1:CA:47:C:H4'	1:CA:48:C:O5'	2.18	0.43
1:CA:501:C:O2'	1:CA:502:A:H5'	2.19	0.43
1:CA:631:C:H5'	1:CA:632:U:O4'	2.19	0.43
1:CA:696:A:H2'	1:CA:697:U:C6	2.54	0.43
1:CA:767:A:H2'	1:CA:768:A:C8	2.54	0.43
18:CB:44:LYS:C	18:CB:47:PRO:HD2	2.39	0.43
3:CD:176:LYS:HD3	3:CD:176:LYS:H	1.83	0.43
8:CI:59:LYS:HZ3	8:CI:60:LEU:CG	2.31	0.43
2:CC:32:LEU:HD21	21:CN:92:ILE:HG12	2.00	0.43
15:CR:55:ALA:O	15:CR:58:ILE:HB	2.18	0.43
32:D4:19:ARG:O	32:D4:21:GLY:N	2.51	0.43
23:DB:1085:A:H1'	23:DB:1105:U:H1'	1.99	0.43
23:DB:175:G:O2'	23:DB:176:A:H5'	2.18	0.43
23:DB:301:G:O5'	46:DU:81:ARG:NH1	2.51	0.43
23:DB:572:A:H3'	23:DB:573:U:O4'	2.18	0.43
23:DB:795:C:H2'	23:DB:796:C:C6	2.54	0.43
26:DD:51:THR:HG22	26:DD:52:THR:N	2.32	0.43
29:DE:170:ARG:NH2	29:DE:176:ASP:HB2	2.33	0.43
29:DE:48:THR:C	29:DE:50:ALA:N	2.71	0.43
29:DE:62:GLN:CG	29:DE:63:LYS:N	2.81	0.43
48:DG:132:LEU:HD13	48:DG:143:VAL:HG23	2.00	0.43
48:DG:5:LYS:HE3	48:DG:61:TRP:CZ3	2.54	0.43
48:DG:64:ALA:O	48:DG:67:ALA:HB3	2.19	0.43
23:DB:1098:A:C2'	24:DI:4:VAL:C	2.87	0.43
24:DI:72:THR:OG1	24:DI:73:PRO:HD2	2.18	0.43
41:DJ:34:ARG:HH11	41:DJ:39:LYS:HG2	1.84	0.43
23:DB:1009:A:OP1	41:DJ:39:LYS:NZ	2.52	0.43
41:DJ:3:THR:HB	41:DJ:44:TYR:HE1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DJ:64:VAL:HG22	41:DJ:68:LYS:HB2	2.01	0.43
27:DK:10:VAL:HG21	27:DK:16:ALA:CA	2.48	0.43
27:DK:36:GLY:HA2	27:DK:62:VAL:O	2.19	0.43
37:DL:110:VAL:HB	37:DL:127:VAL:CG2	2.49	0.43
42:DN:28:LEU:HD23	42:DN:34:ILE:HG13	2.00	0.43
28:DP:26:GLU:HA	28:DP:43:GLU:HA	2.01	0.43
44:DQ:31:TYR:O	44:DQ:32:ARG:C	2.56	0.43
44:DQ:82:LEU:HA	44:DQ:85:ALA:HB3	2.00	0.43
45:DS:20:VAL:C	45:DS:22:ASP:N	2.70	0.43
50:DT:79:ASP:N	50:DT:79:ASP:OD2	2.51	0.43
39:DX:6:LEU:HD22	39:DX:6:LEU:N	2.34	0.43
1:AA:1020:G:N3	1:AA:1020:G:H2'	2.34	0.43
1:AA:1032:G:H2'	1:AA:1033:G:O4'	2.18	0.43
1:AA:1294:G:O2'	1:AA:1295:U:H5'	2.19	0.43
1:AA:1449:C:H2'	1:AA:1450:U:O4'	2.19	0.43
1:AA:238:A:C2'	1:AA:239:U:H5''	2.42	0.43
1:AA:407:U:H2'	1:AA:408:A:C8	2.54	0.43
1:AA:509:A:N3	1:AA:543:U:O2'	2.51	0.43
1:AA:520:A:C2	1:AA:536:C:H1'	2.54	0.43
18:AB:115:ASP:OD1	18:AB:116:LEU:N	2.52	0.43
18:AB:63:LYS:HA	18:AB:224:ARG:HH12	1.84	0.43
2:AC:51:VAL:HG23	2:AC:68:HIS:O	2.18	0.43
3:AD:107:GLY:HA2	3:AD:112:GLU:OE2	2.17	0.43
3:AD:154:VAL:O	3:AD:158:LEU:HG	2.19	0.43
3:AD:187:ARG:HH11	3:AD:191:SER:HA	1.84	0.43
4:AE:156:ARG:HB3	7:AH:43:GLY:O	2.19	0.43
4:AE:43:GLY:O	4:AE:72:ASN:HA	2.18	0.43
5:AF:10:VAL:HA	5:AF:84:VAL:HA	2.00	0.43
8:AI:12:LYS:NZ	8:AI:12:LYS:HB2	2.34	0.43
8:AI:55:ASP:HB2	8:AI:59:LYS:CD	2.42	0.43
10:AK:95:THR:HG23	10:AK:96:ILE:HG13	2.01	0.43
12:AM:21:ILE:HG22	12:AM:23:GLY:N	2.29	0.43
12:AM:79:LEU:HA	12:AM:82:LEU:CD1	2.47	0.43
19:AU:8:ASN:O	19:AU:9:GLU:HB3	2.19	0.43
34:B3:21:PHE:CE1	34:B3:58:ILE:HG23	2.53	0.43
23:BB:1259:G:H2'	23:BB:1260:A:H8	1.82	0.43
23:BB:1850:G:H2'	23:BB:1851:U:C6	2.53	0.43
23:BB:202:U:H2'	23:BB:203:A:O4'	2.18	0.43
23:BB:2770:G:H8	23:BB:2770:G:O5'	2.02	0.43
23:BB:2853:C:H2'	23:BB:2854:G:C8	2.53	0.43
23:BB:528:A:H2'	23:BB:529:A:H5''	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:572:A:H3'	23:BB:573:U:O4'	2.18	0.43
23:BB:1819:A:OP1	25:BC:159:THR:HG21	2.18	0.43
25:BC:106:PRO:HA	25:BC:195:GLY:N	2.33	0.43
29:BE:161:ALA:HA	29:BE:164:LEU:HB2	2.01	0.43
29:BE:62:GLN:HB2	29:BE:62:GLN:HE21	1.52	0.43
23:BB:659:G:H4'	29:BE:95:LYS:HB3	2.00	0.43
47:BF:146:ASP:O	47:BF:147:ARG:C	2.56	0.43
47:BF:64:PRO:HA	47:BF:88:VAL:HG22	2.00	0.43
48:BG:54:ARG:HA	48:BG:54:ARG:HE	1.83	0.43
48:BG:93:TYR:CD2	48:BG:93:TYR:N	2.86	0.43
23:BB:1061:U:O4	24:BI:10:LEU:HA	2.18	0.43
41:BJ:12:LYS:O	41:BJ:13:ARG:HB2	2.19	0.43
27:BK:10:VAL:HG21	27:BK:16:ALA:CA	2.49	0.43
38:BM:19:GLY:N	38:BM:38:ARG:NH1	2.66	0.43
43:BO:105:ALA:C	43:BO:107:ALA:H	2.19	0.43
43:BO:7:ARG:N	43:BO:10:ARG:HH11	2.17	0.43
49:BR:15:SER:H	49:BR:18:GLN:CG	2.32	0.43
45:BS:52:GLU:HA	45:BS:55:ILE:HG22	2.01	0.43
45:BS:86:MET:HB3	45:BS:94:ASP:HB2	2.00	0.43
23:BB:856:G:C1'	52:BW:23:LYS:HB3	2.48	0.43
23:BB:200:U:O2'	51:BZ:22:LEU:HD12	2.19	0.43
51:BZ:33:LEU:HA	51:BZ:51:VAL:O	2.19	0.43
1:CA:1180:A:OP1	8:CI:104:THR:HG22	2.19	0.43
1:CA:1347:G:H5''	8:CI:109:GLN:O	2.18	0.43
1:CA:332:G:H2'	1:CA:333:U:H6	1.84	0.43
1:CA:410:G:P	3:CD:25:ARG:HD2	2.58	0.43
1:CA:484:G:H3'	1:CA:484:G:OP2	2.17	0.43
1:CA:502:A:H2'	1:CA:503:C:O4'	2.19	0.43
1:CA:947:G:H4'	1:CA:1332:A:H2	1.83	0.43
1:CA:950:U:H2'	1:CA:951:G:H8	1.83	0.43
1:CA:955:U:H2'	1:CA:956:U:O4'	2.18	0.43
18:CB:169:HIS:N	18:CB:169:HIS:CD2	2.85	0.43
2:CC:129:PHE:CZ	2:CC:130:ARG:HD2	2.54	0.43
3:CD:155:LYS:HG2	3:CD:156:ALA:H	1.83	0.43
3:CD:187:ARG:O	3:CD:191:SER:N	2.52	0.43
4:CE:141:ASP:HA	4:CE:144:GLU:OE1	2.19	0.43
10:CK:67:GLU:HG3	10:CK:68:ARG:N	2.33	0.43
11:CL:106:VAL:CG2	11:CL:116:TYR:HB3	2.48	0.43
12:CM:22:TYR:HD1	12:CM:65:GLU:HA	1.82	0.43
12:CM:85:TYR:HB2	16:CS:72:GLU:HB2	2.00	0.43
21:CN:60:ARG:NH2	21:CN:69:PRO:HB3	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CO:16:GLY:HA2	20:CO:27:VAL:CG2	2.48	0.43
20:CO:7:ALA:O	20:CO:11:ILE:HG23	2.19	0.43
20:CO:64:ARG:NH2	20:CO:88:ARG:HH11	2.17	0.43
14:CQ:26:ARG:N	14:CQ:39:ARG:O	2.52	0.43
14:CQ:58:VAL:HG12	14:CQ:77:VAL:HG13	2.00	0.43
1:CA:719:C:H2'	15:CR:38:ILE:CD1	2.49	0.43
16:CS:29:PRO:HB3	16:CS:47:THR:HG22	2.00	0.43
32:D4:8:LYS:HG2	32:D4:9:LYS:H	1.83	0.43
22:DA:109:A:H2'	22:DA:110:C:H6	1.84	0.43
22:DA:111:U:H2'	22:DA:112:G:C8	2.53	0.43
22:DA:25:U:H4'	22:DA:26:C:OP1	2.17	0.43
22:DA:67:G:H2'	22:DA:68:C:H6	1.84	0.43
22:DA:6:G:O2'	22:DA:7:G:H5'	2.19	0.43
23:DB:1062:G:H2'	23:DB:1063:G:H8	1.83	0.43
23:DB:1463:C:H2'	23:DB:1464:G:H8	1.83	0.43
23:DB:1486:U:H2'	23:DB:1487:U:H6	1.83	0.43
23:DB:1541:C:O2'	23:DB:1542:U:H5'	2.19	0.43
23:DB:1600:C:C2'	23:DB:1601:G:H5'	2.48	0.43
23:DB:163:C:H2'	23:DB:164:C:H6	1.84	0.43
23:DB:195:A:H5''	37:DL:47:ARG:NH2	2.34	0.43
23:DB:2376:A:C2	43:DO:92:PHE:HB3	2.54	0.43
23:DB:286:U:H2'	23:DB:287:G:C8	2.51	0.43
23:DB:2869:G:H2'	23:DB:2870:C:O4'	2.19	0.43
23:DB:690:G:H2'	23:DB:691:C:O4'	2.18	0.43
23:DB:744:U:H2'	23:DB:745:G:C8	2.54	0.43
23:DB:735:A:N7	23:DB:761:A:H2	2.16	0.43
26:DD:124:ARG:HE	26:DD:125:TRP:HE1	1.64	0.43
29:DE:1:MET:O	29:DE:13:THR:HA	2.18	0.43
48:DG:93:TYR:O	48:DG:94:ARG:O	2.36	0.43
24:DI:4:VAL:O	24:DI:5:GLN:O	2.36	0.43
24:DI:54:ILE:HD11	24:DI:71:LYS:N	2.33	0.43
23:DB:958:U:O4	38:DM:16:ARG:HA	2.18	0.43
38:DM:59:ARG:HE	38:DM:60:GLN:CG	2.11	0.43
43:DO:66:GLY:CA	43:DO:102:ARG:HH21	2.17	0.43
28:DP:58:PHE:HD1	28:DP:73:PHE:HD2	1.65	0.43
49:DR:49:ILE:HG21	49:DR:54:VAL:HA	2.01	0.43
45:DS:60:HIS:ND1	45:DS:60:HIS:O	2.51	0.43
45:DS:73:LYS:HD2	45:DS:73:LYS:HA	1.66	0.43
46:DU:41:VAL:O	46:DU:59:GLU:HG3	2.17	0.43
35:DV:68:LYS:HG2	35:DV:69:GLU:N	2.33	0.43
1:AA:1308:U:H3'	12:AM:97:ARG:HH12	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1509:C:H2'	1:AA:1510:C:H6	1.84	0.43
1:AA:123:U:OP1	1:AA:312:C:H5'	2.18	0.43
18:AB:45:THR:HG22	18:AB:49:PHE:CZ	2.54	0.43
18:AB:55:GLU:HG3	18:AB:197:PHE:CE1	2.54	0.43
3:AD:96:ARG:O	3:AD:100:VAL:HG23	2.18	0.43
6:AG:43:TYR:HA	6:AG:46:LEU:HB3	2.00	0.43
7:AH:9:MET:O	7:AH:13:ILE:HG13	2.19	0.43
9:AJ:18:ILE:HG23	9:AJ:19:ASP:N	2.33	0.43
21:AN:41:TRP:HB3	21:AN:44:VAL:CG2	2.49	0.43
20:AO:45:GLU:O	20:AO:47:LYS:N	2.47	0.43
16:AS:35:ARG:HB2	16:AS:71:GLY:CA	2.44	0.43
16:AS:38:THR:HA	16:AS:69:LYS:HD3	2.00	0.43
22:BA:75:G:H2'	22:BA:76:G:C8	2.50	0.43
23:BB:1197:G:H5'	23:BB:1227:G:O2'	2.19	0.43
23:BB:1252:G:H1'	44:BQ:32:ARG:NH2	2.30	0.43
23:BB:1535:A:O2'	23:BB:1537:G:N7	2.49	0.43
23:BB:163:C:O4'	23:BB:163:C:O2	2.36	0.43
23:BB:1662:U:O2	23:BB:2687:U:H4'	2.19	0.43
23:BB:1751:U:H2'	23:BB:1752:C:C6	2.53	0.43
23:BB:182:A:H2'	23:BB:183:C:C6	2.53	0.43
23:BB:1860:G:O2'	23:BB:1861:G:H5'	2.19	0.43
23:BB:225:C:H2'	23:BB:226:A:O4'	2.18	0.43
23:BB:222:A:H61	23:BB:232:G:H1'	1.84	0.43
23:BB:233:A:H61	23:BB:428:A:N6	2.16	0.43
23:BB:2379:G:H2'	23:BB:2380:C:C6	2.53	0.43
23:BB:2437:G:H2'	23:BB:2438:U:C6	2.54	0.43
23:BB:2895:G:O2'	23:BB:2896:C:H5'	2.18	0.43
23:BB:644:A:O2'	23:BB:645:C:H2'	2.18	0.43
23:BB:816:C:O2'	23:BB:817:C:H5'	2.19	0.43
23:BB:850:U:O2'	30:BY:22:THR:HG22	2.18	0.43
23:BB:988:A:O5'	30:BY:11:SER:HB3	2.19	0.43
26:BD:13:ARG:HH12	28:BP:74:GLN:NE2	2.17	0.43
29:BE:114:ARG:HG3	29:BE:114:ARG:NH1	2.33	0.43
29:BE:148:ILE:HD13	29:BE:187:VAL:HG23	2.00	0.43
22:BA:54:G:N2	47:BF:25:MET:HG2	2.31	0.43
40:BH:90:LEU:HD11	40:BH:146:VAL:CG1	2.49	0.43
24:BI:49:GLU:HG3	24:BI:54:ILE:HD11	2.01	0.43
41:BJ:88:THR:O	41:BJ:91:GLU:HB2	2.18	0.43
27:BK:62:VAL:HG12	27:BK:63:VAL:N	2.34	0.43
23:BB:1276:A:H1'	42:BN:16:HIS:HE2	1.83	0.43
42:BN:31:HIS:O	42:BN:32:GLU:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BP:3:ILE:HD13	28:BP:3:ILE:C	2.39	0.43
28:BP:77:SER:O	28:BP:80:VAL:HG12	2.19	0.43
28:BP:95:LYS:HG3	28:BP:97:TYR:CE1	2.52	0.43
49:BR:82:HIS:O	49:BR:83:TYR:C	2.58	0.43
50:BT:87:LEU:O	50:BT:88:LYS:O	2.37	0.43
35:BV:87:GLN:HE21	35:BV:87:GLN:HB2	1.65	0.43
52:BW:25:PHE:C	52:BW:27:GLY:N	2.72	0.43
52:BW:32:ALA:C	52:BW:34:SER:N	2.71	0.43
23:BB:77:G:H5'	39:BX:52:ARG:HG2	2.00	0.43
39:BX:49:ASP:O	39:BX:53:VAL:HG23	2.18	0.43
1:CA:1103:C:H2'	1:CA:1104:G:O4'	2.18	0.43
1:CA:1109:C:H2'	1:CA:1110:A:O4'	2.18	0.43
1:CA:1115:U:H6	1:CA:1115:U:O5'	2.01	0.43
1:CA:1382:C:O2'	1:CA:1383:C:H5'	2.18	0.43
1:CA:1426:G:H2'	1:CA:1427:C:O4'	2.19	0.43
1:CA:197:A:H4'	1:CA:198:G:O5'	2.17	0.43
1:CA:213:G:C8	1:CA:214:C:C5	3.06	0.43
1:CA:521:G:O2'	1:CA:522:C:H5'	2.19	0.43
1:CA:537:G:H2'	1:CA:538:G:H8	1.83	0.43
1:CA:708:C:O2'	1:CA:709:U:H5'	2.19	0.43
2:CC:110:LEU:HD22	2:CC:145:ALA:HB2	2.00	0.43
2:CC:183:TYR:O	2:CC:184:ASN:HB2	2.17	0.43
3:CD:108:ALA:H	3:CD:112:GLU:CD	2.22	0.43
3:CD:117:VAL:HA	3:CD:122:ILE:HG12	2.01	0.43
3:CD:181:PHE:O	3:CD:182:LYS:C	2.57	0.43
6:CG:43:TYR:O	6:CG:44:SER:C	2.56	0.43
9:CJ:7:ARG:HG3	9:CJ:102:LEU:O	2.17	0.43
9:CJ:10:LEU:HB2	9:CJ:72:ARG:HB2	2.01	0.43
1:CA:1254:A:OP1	9:CJ:47:GLU:HG3	2.19	0.43
10:CK:86:LYS:HB2	10:CK:113:THR:HA	2.00	0.43
20:CO:70:LEU:HD11	20:CO:77:ARG:CB	2.45	0.43
1:CA:564:C:N1	14:CQ:32:ILE:HD11	2.34	0.43
14:CQ:40:THR:HG22	14:CQ:41:THR:N	2.34	0.43
14:CQ:61:ARG:O	14:CQ:61:ARG:HG3	2.19	0.43
15:CR:27:THR:O	15:CR:28:LEU:HD23	2.18	0.43
16:CS:33:TRP:CD1	16:CS:51:HIS:HB3	2.53	0.43
33:D1:8:ILE:HD12	33:D1:51:ALA:HA	1.98	0.43
22:DA:52:A:C2'	22:DA:53:A:H5'	2.48	0.43
23:DB:121:G:H2'	23:DB:122:G:C8	2.53	0.43
23:DB:1241:A:N3	23:DB:1241:A:H5'	2.34	0.43
23:DB:1328:A:H2'	23:DB:1330:C:C5	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1464:G:H2'	23:DB:1465:G:C8	2.53	0.43
23:DB:1784:A:H4'	23:DB:1785:A:O5'	2.18	0.43
23:DB:1841:U:C2	23:DB:1842:G:C8	3.07	0.43
23:DB:2011:U:O2'	23:DB:2012:G:H5'	2.19	0.43
23:DB:210:C:O2'	23:DB:211:C:H5'	2.18	0.43
23:DB:2785:C:H2'	23:DB:2786:U:O4'	2.19	0.43
23:DB:2806:C:H2'	23:DB:2807:U:O4'	2.19	0.43
23:DB:2815:C:H2'	23:DB:2816:G:H8	1.83	0.43
23:DB:677:A:H2'	23:DB:678:C:C6	2.52	0.43
23:DB:729:G:H5''	23:DB:730:A:H5''	2.00	0.43
25:DC:124:LYS:HG3	25:DC:125:PRO:HD2	2.00	0.43
26:DD:40:LEU:HD12	26:DD:41:ALA:N	2.34	0.43
29:DE:148:ILE:HD13	29:DE:187:VAL:HG23	1.99	0.43
29:DE:24:ASN:OD1	29:DE:27:LEU:HB2	2.18	0.43
29:DE:58:LYS:N	29:DE:58:LYS:NZ	2.57	0.43
29:DE:67:ARG:NH1	29:DE:70:SER:OG	2.51	0.43
29:DE:40:ARG:NH2	29:DE:92:HIS:NE2	2.66	0.43
47:DF:137:PHE:CD2	47:DF:137:PHE:N	2.87	0.43
40:DH:99:ILE:O	40:DH:103:VAL:HG12	2.19	0.43
24:DI:68:PHE:CD1	24:DI:68:PHE:N	2.87	0.43
41:DJ:40:HIS:O	44:DQ:66:ALA:HB1	2.18	0.43
41:DJ:4:PHE:HB3	41:DJ:44:TYR:CE1	2.53	0.43
37:DL:111:ILE:HG22	37:DL:112:LEU:N	2.31	0.43
42:DN:31:HIS:N	42:DN:31:HIS:ND1	2.67	0.43
44:DQ:91:ARG:HE	44:DQ:93:ILE:CG2	2.32	0.43
49:DR:69:GLY:O	49:DR:90:ARG:HG3	2.19	0.43
50:DT:22:THR:HA	50:DT:25:GLU:HB3	2.01	0.43
46:DU:18:LYS:HE2	46:DU:18:LYS:C	2.39	0.43
52:DW:58:LEU:N	52:DW:58:LEU:HD22	2.34	0.43
30:DY:40:THR:O	30:DY:43:ILE:HG23	2.19	0.43
1:AA:1164:G:O2'	1:AA:1165:U:H5'	2.18	0.43
1:AA:1201:A:H1'	1:AA:1202:U:OP2	2.19	0.43
1:AA:1292:G:H2'	1:AA:1293:C:H6	1.84	0.43
1:AA:1316:G:C2	1:AA:1318:A:H5''	2.53	0.43
1:AA:1323:G:H2'	1:AA:1324:A:H8	1.81	0.43
1:AA:1342:C:O2'	1:AA:1343:G:H5'	2.18	0.43
1:AA:1509:C:O2'	1:AA:1510:C:H5'	2.19	0.43
1:AA:229:U:H2'	1:AA:230:G:H8	1.82	0.43
1:AA:920:U:O2'	1:AA:921:U:H5'	2.18	0.43
18:AB:101:THR:HG22	18:AB:174:GLU:OE1	2.18	0.43
18:AB:19:THR:OG1	18:AB:20:ARG:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:67:LEU:HD21	18:AB:91:VAL:HG23	2.00	0.43
2:AC:27:GLU:HB2	2:AC:31:ASN:OD1	2.19	0.43
2:AC:54:ILE:O	2:AC:54:ILE:HG12	2.19	0.43
3:AD:157:ALA:C	3:AD:159:GLU:N	2.71	0.43
7:AH:23:ALA:CB	7:AH:61:THR:HA	2.49	0.43
7:AH:25:THR:HG22	7:AH:26:MET:H	1.84	0.43
8:AI:119:LYS:HD2	8:AI:119:LYS:O	2.19	0.43
8:AI:11:ARG:HE	8:AI:12:LYS:CD	2.31	0.43
8:AI:51:LEU:CB	8:AI:56:MET:SD	2.99	0.43
8:AI:96:GLU:HA	8:AI:99:LYS:HE3	2.00	0.43
9:AJ:57:VAL:HG13	9:AJ:58:ASN:N	2.32	0.43
1:AA:707:U:H4'	10:AK:21:HIS:CG	2.53	0.43
10:AK:91:GLY:C	10:AK:93:GLU:N	2.72	0.43
11:AL:24:GLU:C	11:AL:26:CYS:H	2.19	0.43
12:AM:69:ARG:HA	12:AM:72:ILE:HD12	2.00	0.43
14:AQ:34:GLY:O	14:AQ:35:LYS:C	2.57	0.43
34:B3:61:LEU:N	34:B3:62:PRO:CD	2.82	0.43
22:BA:76:G:H1	22:BA:101:A:N6	2.16	0.43
23:BB:1570:A:H2'	23:BB:1571:A:C8	2.54	0.43
23:BB:1727:C:H2'	23:BB:1728:C:H6	1.82	0.43
23:BB:1989:G:H2'	23:BB:1990:C:O4'	2.19	0.43
23:BB:2626:C:O2'	23:BB:2627:G:H5'	2.19	0.43
23:BB:2721:A:H2'	23:BB:2722:G:H8	1.82	0.43
23:BB:441:U:O2'	23:BB:442:G:H5'	2.18	0.43
23:BB:479:A:H1'	23:BB:480:A:H5''	2.00	0.43
23:BB:481:G:C2	23:BB:507:A:C4	3.07	0.43
23:BB:559:G:H1'	44:BQ:55:GLN:NE2	2.34	0.43
23:BB:765:C:O2'	23:BB:766:U:H5'	2.18	0.43
23:BB:768:G:O2'	23:BB:769:U:H5'	2.18	0.43
26:BD:2:ILE:HG23	26:BD:84:LEU:HB3	2.01	0.43
24:BI:19:PRO:HG2	24:BI:22:PRO:HB2	2.01	0.43
41:BJ:77:HIS:HD2	41:BJ:84:ILE:HB	1.83	0.43
27:BK:112:PHE:O	27:BK:113:MET:C	2.57	0.43
37:BL:78:ARG:HB3	37:BL:113:ALA:CB	2.38	0.43
42:BN:116:VAL:HG13	42:BN:117:ASP:N	2.33	0.43
42:BN:118:ARG:HE	42:BN:118:ARG:HB3	1.49	0.43
42:BN:29:VAL:HG13	42:BN:83:LEU:CD2	2.49	0.43
43:BO:7:ARG:HG3	43:BO:96:GLY:O	2.19	0.43
44:BQ:30:VAL:HG11	44:BQ:33:VAL:HG22	2.00	0.43
35:BV:68:LYS:HG2	35:BV:69:GLU:N	2.33	0.43
52:BW:8:SER:O	52:BW:9:THR:HB	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1057:G:H2'	1:CA:1058:G:C8	2.54	0.43
1:CA:1016:A:H4'	1:CA:1218:C:C4'	2.49	0.43
1:CA:1253:G:OP1	9:CJ:46:LYS:HG2	2.19	0.43
1:CA:222:C:H2'	1:CA:223:A:C8	2.54	0.43
1:CA:634:C:H6	1:CA:634:C:O5'	2.02	0.43
1:CA:771:G:H2'	1:CA:772:U:C6	2.54	0.43
18:CB:114:LYS:HZ1	18:CB:151:LYS:HG3	1.84	0.43
3:CD:187:ARG:HH11	3:CD:191:SER:HA	1.83	0.43
1:CA:16:A:O2'	4:CE:20:VAL:HG22	2.19	0.43
4:CE:80:LEU:HG	4:CE:122:VAL:HG11	2.01	0.43
6:CG:46:LEU:HG	6:CG:57:GLU:HB3	2.00	0.43
9:CJ:10:LEU:O	9:CJ:71:LEU:HA	2.19	0.43
11:CL:54:VAL:HG12	11:CL:55:ARG:N	2.33	0.43
12:CM:52:ILE:HG23	12:CM:53:ASP:H	1.83	0.43
13:CP:36:VAL:O	13:CP:36:VAL:HG13	2.18	0.43
22:DA:9:G:O2'	22:DA:10:G:H5'	2.18	0.43
23:DB:52:A:C5	23:DB:118:A:C2	3.06	0.43
23:DB:1360:G:C2'	23:DB:1361:G:H5'	2.48	0.43
23:DB:1400:U:O2'	23:DB:1401:G:H5'	2.19	0.43
23:DB:1468:U:HO2'	23:DB:1469:A:H8	1.62	0.43
23:DB:1310:G:N2	23:DB:1610:A:H8	2.04	0.43
23:DB:1726:C:H2'	23:DB:1727:C:C6	2.54	0.43
23:DB:1727:C:H2'	23:DB:1728:C:O4'	2.19	0.43
23:DB:2348:U:H4'	33:D1:40:PRO:HG2	2.00	0.43
23:DB:2709:G:H2'	23:DB:2710:C:C6	2.53	0.43
23:DB:2628:C:O2'	23:DB:2781:A:H2'	2.18	0.43
23:DB:2852:G:O2'	23:DB:2853:C:H5'	2.19	0.43
23:DB:682:G:O2'	23:DB:683:U:H5'	2.18	0.43
23:DB:1824:G:OP2	25:DC:52:HIS:CE1	2.72	0.43
25:DC:93:VAL:CG1	25:DC:94:LEU:H	2.17	0.43
29:DE:148:ILE:HD13	29:DE:187:VAL:CG2	2.49	0.43
47:DF:11:VAL:O	47:DF:13:LYS:N	2.52	0.43
40:DH:26:ALA:O	40:DH:28:ASN:N	2.52	0.43
41:DJ:83:GLY:O	41:DJ:84:ILE:C	2.56	0.43
23:DB:1190:G:P	37:DL:32:GLY:HA2	2.58	0.43
37:DL:91:ASP:HB2	37:DL:94:THR:HG23	2.01	0.43
38:DM:130:PHE:O	38:DM:131:VAL:HG13	2.19	0.43
38:DM:73:ILE:HG21	38:DM:91:TYR:OH	2.19	0.43
28:DP:6:GLN:HE21	28:DP:7:LEU:HD12	1.84	0.43
44:DQ:104:ALA:HA	49:DR:46:GLU:CD	2.38	0.43
44:DQ:57:ARG:C	44:DQ:59:LEU:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DS:20:VAL:O	45:DS:22:ASP:N	2.52	0.43
45:DS:43:ALA:O	45:DS:47:VAL:HG23	2.18	0.43
46:DU:1:ALA:HB1	46:DU:84:PHE:CZ	2.54	0.43
39:DX:17:GLU:CD	39:DX:18:LEU:N	2.72	0.43
1:AA:1084:G:H2'	1:AA:1085:U:C6	2.54	0.43
1:AA:1096:C:O2'	1:AA:1097:C:H5'	2.19	0.43
1:AA:1190:G:OP1	2:AC:4:VAL:N	2.44	0.43
1:AA:415:A:O4'	1:AA:415:A:N3	2.52	0.43
1:AA:656:G:H2'	1:AA:657:U:C6	2.54	0.43
1:AA:78:A:O2'	1:AA:79:G:H5'	2.19	0.43
1:AA:920:U:C2	1:AA:921:U:C5	3.07	0.43
18:AB:100:LEU:HD21	18:AB:160:LEU:HD12	2.00	0.43
2:AC:152:VAL:HG13	2:AC:197:VAL:CG2	2.49	0.43
4:AE:33:THR:HB	4:AE:49:TYR:OH	2.19	0.43
9:AJ:9:ARG:HB2	9:AJ:99:GLN:HB3	2.01	0.43
12:AM:56:ARG:O	12:AM:59:VAL:HG12	2.19	0.43
2:AC:11:LEU:HD12	21:AN:96:LYS:HA	2.01	0.43
22:BA:64:G:H2'	22:BA:65:U:H6	1.82	0.43
23:BB:1057:A:H62	23:BB:1086:A:H2'	1.84	0.43
23:BB:1135:C:N4	23:BB:1139:G:C6	2.87	0.43
23:BB:1326:U:H2'	23:BB:1327:A:O4'	2.19	0.43
23:BB:1486:U:O2'	23:BB:1487:U:H5'	2.19	0.43
23:BB:2267:A:OP2	23:BB:2268:A:H5''	2.18	0.43
23:BB:2272:U:O2'	23:BB:2273:A:H8	2.02	0.43
23:BB:245:G:O2'	23:BB:246:C:H5'	2.19	0.43
23:BB:2602:A:H3'	23:BB:2602:A:OP1	2.19	0.43
23:BB:2604:U:O2'	23:BB:2605:U:H5'	2.19	0.43
23:BB:325:G:H2'	23:BB:326:G:H8	1.83	0.43
23:BB:311:A:H2	23:BB:331:C:H5''	1.83	0.43
23:BB:666:A:H4'	37:BL:48:ARG:HD2	2.01	0.43
23:BB:704:G:H1'	23:BB:727:A:N6	2.33	0.43
23:BB:950:G:H2'	23:BB:951:C:H6	1.79	0.43
23:BB:969:G:H2'	23:BB:970:U:H6	1.83	0.43
26:BD:148:GLN:HB2	26:BD:149:ASN:H	1.62	0.43
26:BD:40:LEU:HD12	26:BD:41:ALA:N	2.34	0.43
29:BE:105:LEU:HD21	29:BE:177:PRO:HB3	2.00	0.43
47:BF:111:ARG:H	47:BF:111:ARG:HD2	1.84	0.43
40:BH:84:ALA:HA	40:BH:90:LEU:CA	2.29	0.43
38:BM:21:ALA:CB	38:BM:100:LYS:HG2	2.49	0.43
38:BM:111:GLU:HA	38:BM:114:ARG:NH2	2.34	0.43
38:BM:126:ILE:HG22	38:BM:127:LYS:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BP:19:PHE:CE2	28:BP:83:ILE:HD11	2.54	0.43
44:BQ:49:ARG:O	44:BQ:53:LYS:HE2	2.19	0.43
23:BB:584:C:P	44:BQ:5:ARG:HD3	2.58	0.43
50:BT:21:SER:O	50:BT:25:GLU:HB2	2.19	0.43
46:BU:41:VAL:O	46:BU:59:GLU:HG3	2.19	0.43
35:BV:75:GLN:CB	35:BV:90:ASP:HB2	2.49	0.43
51:BZ:49:LEU:O	51:BZ:51:VAL:HG23	2.19	0.43
1:CA:1023:U:C2'	1:CA:1024:G:H5'	2.49	0.43
1:CA:1190:G:OP2	2:CC:4:VAL:HB	2.18	0.43
1:CA:1298:U:H1'	1:CA:1299:A:C2	2.53	0.43
1:CA:1298:U:H1'	1:CA:1299:A:N1	2.34	0.43
1:CA:1374:A:O3'	6:CG:27:ASN:ND2	2.51	0.43
3:CD:107:GLY:C	3:CD:157:ALA:HB1	2.38	0.43
3:CD:49:ASP:OD2	4:CE:111:ARG:HD2	2.18	0.43
7:CH:9:MET:O	7:CH:13:ILE:HG13	2.19	0.43
8:CI:35:GLU:HA	8:CI:39:GLY:CA	2.49	0.43
9:CJ:16:ARG:H	9:CJ:16:ARG:HG3	1.52	0.43
9:CJ:42:LEU:HB3	9:CJ:71:LEU:HD11	2.01	0.43
11:CL:27:PRO:O	11:CL:28:GLN:NE2	2.52	0.43
12:CM:28:ARG:O	12:CM:32:ILE:HB	2.19	0.43
21:CN:3:GLN:HA	21:CN:6:LYS:HG3	2.00	0.43
20:CO:14:GLU:CB	20:CO:84:ARG:HH22	2.32	0.43
16:CS:16:LYS:HD2	16:CS:17:LYS:HZ2	1.82	0.43
33:D1:10:LEU:HD23	33:D1:35:LEU:HD21	2.00	0.43
33:D1:26:LYS:HG2	33:D1:52:LYS:HD2	2.01	0.43
36:D2:26:ASN:O	36:D2:30:VAL:HG23	2.18	0.43
34:D3:57:VAL:O	34:D3:59:ALA:N	2.47	0.43
22:DA:64:G:O2'	22:DA:65:U:H5'	2.19	0.43
23:DB:1099:G:C5'	24:DI:4:VAL:HG12	2.48	0.43
23:DB:1459:G:H8	23:DB:1461:C:H5	1.67	0.43
23:DB:2282:G:O2'	23:DB:2283:C:OP2	2.33	0.43
23:DB:2336:A:O2'	23:DB:2337:G:P	2.77	0.43
23:DB:2494:G:H2'	23:DB:2495:G:H8	1.83	0.43
23:DB:2552:U:H3'	23:DB:2554:U:OP2	2.19	0.43
23:DB:322:A:H1'	23:DB:339:U:O2	2.19	0.43
23:DB:479:A:O2'	23:DB:481:G:H5'	2.18	0.43
23:DB:68:G:H2'	23:DB:69:C:C6	2.54	0.43
23:DB:707:G:O2'	23:DB:708:G:H5'	2.18	0.43
23:DB:1568:G:H4'	25:DC:58:LYS:HD3	2.00	0.43
26:DD:37:VAL:HG12	26:DD:38:LYS:N	2.33	0.43
29:DE:24:ASN:HD22	29:DE:24:ASN:C	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:99:PHE:O	47:DF:102:LEU:HB3	2.18	0.43
47:DF:64:PRO:HA	47:DF:88:VAL:HG22	2.00	0.43
40:DH:128:HIS:N	40:DH:144:VAL:O	2.52	0.43
40:DH:114:GLU:CG	40:DH:133:GLN:HB2	2.49	0.43
40:DH:2:GLN:HB2	40:DH:39:ALA:CB	2.49	0.43
24:DI:101:SER:OG	24:DI:104:GLN:HG3	2.19	0.43
24:DI:96:LYS:HD3	24:DI:138:VAL:HG21	2.00	0.43
41:DJ:133:ALA:C	41:DJ:135:GLN:H	2.22	0.43
41:DJ:12:LYS:O	41:DJ:13:ARG:HB2	2.19	0.43
41:DJ:13:ARG:HB3	41:DJ:53:TYR:HD2	1.83	0.43
41:DJ:15:TRP:CD2	41:DJ:53:TYR:HB3	2.54	0.43
27:DK:54:LYS:HE3	27:DK:54:LYS:HB3	1.87	0.43
37:DL:127:VAL:HG22	37:DL:131:ALA:HB3	2.00	0.43
42:DN:11:ASN:HA	42:DN:11:ASN:HD22	1.52	0.43
42:DN:36:THR:OG1	42:DN:40:LYS:HD2	2.19	0.43
28:DP:85:VAL:HG21	28:DP:88:ARG:NH1	2.33	0.43
49:DR:31:GLU:O	49:DR:63:VAL:HG22	2.19	0.43
49:DR:14:VAL:HG21	49:DR:98:ILE:CD1	2.49	0.43
49:DR:98:ILE:HG22	49:DR:99:THR:N	2.34	0.43
45:DS:33:LEU:HD23	45:DS:51:LEU:HD23	2.00	0.43
45:DS:83:LYS:HB3	45:DS:95:ARG:NH1	2.34	0.43
51:DZ:40:VAL:CG2	51:DZ:43:GLU:HB3	2.44	0.43
1:AA:1117:A:C2	1:AA:1180:A:H1'	2.53	0.42
1:AA:1360:A:N6	1:AA:1361:G:C2	2.87	0.42
1:AA:454:G:H2'	1:AA:455:G:H8	1.83	0.42
1:AA:631:C:H5''	1:AA:632:U:O4'	2.19	0.42
1:AA:861:G:O2'	1:AA:862:C:H5'	2.18	0.42
1:AA:986:U:H1'	16:AS:54:ARG:HA	2.00	0.42
18:AB:83:ALA:O	18:AB:88:GLN:HB2	2.19	0.42
1:AA:403:C:H5'	3:AD:131:ILE:HG23	2.00	0.42
3:AD:165:GLU:CD	3:AD:166:LYS:H	2.22	0.42
3:AD:4:LEU:N	3:AD:4:LEU:HD12	2.33	0.42
1:AA:16:A:O2'	4:AE:20:VAL:HG22	2.19	0.42
5:AF:51:ILE:HD11	5:AF:86:ARG:HG3	2.01	0.42
1:AA:939:G:C5'	6:AG:101:ARG:HH12	2.20	0.42
6:AG:128:GLU:HG2	6:AG:130:LYS:HB2	2.00	0.42
1:AA:878:A:C5'	7:AH:80:PRO:HG2	2.49	0.42
21:AN:27:LYS:C	21:AN:29:ILE:H	2.22	0.42
21:AN:51:PRO:HB2	21:AN:54:SER:HB3	2.01	0.42
20:AO:27:VAL:O	20:AO:30:ALA:HB3	2.19	0.42
13:AP:74:LEU:HA	13:AP:77:GLU:OE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AQ:4:ILE:HD12	14:AQ:4:ILE:O	2.19	0.42
16:AS:43:MET:O	16:AS:46:LEU:HD23	2.18	0.42
10:AK:125:LYS:O	19:AU:33:ARG:CZ	2.67	0.42
33:B1:9:LYS:O	33:B1:50:GLU:HG3	2.18	0.42
23:BB:1005:C:H2'	23:BB:1006:C:H6	1.84	0.42
23:BB:1312:U:O2	23:BB:1603:A:N1	2.52	0.42
23:BB:1317:G:H2'	23:BB:1318:U:O4'	2.19	0.42
23:BB:1457:U:H5'	23:BB:1458:U:OP1	2.19	0.42
23:BB:1572:A:O2'	23:BB:1573:G:H5'	2.19	0.42
23:BB:1660:G:H2'	23:BB:1661:G:H8	1.83	0.42
23:BB:188:G:O2'	23:BB:189:G:H5'	2.19	0.42
23:BB:235:U:H2'	23:BB:236:C:H6	1.83	0.42
23:BB:2665:A:C2'	23:BB:2666:C:H5'	2.49	0.42
23:BB:527:C:C4	23:BB:2779:U:H2'	2.54	0.42
23:BB:247:G:H4'	23:BB:386:G:C4	2.54	0.42
23:BB:603:A:N6	23:BB:655:A:O4'	2.52	0.42
25:BC:157:ALA:C	25:BC:159:THR:H	2.22	0.42
29:BE:146:VAL:HG12	29:BE:147:LEU:N	2.34	0.42
47:BF:40:GLY:HA2	47:BF:84:ILE:O	2.18	0.42
47:BF:89:THR:O	47:BF:91:ARG:CZ	2.66	0.42
24:BI:63:ASP:O	24:BI:65:SER:N	2.52	0.42
37:BL:111:ILE:HG22	37:BL:112:LEU:N	2.33	0.42
37:BL:95:LEU:H	37:BL:95:LEU:HG	1.60	0.42
28:BP:26:GLU:HA	28:BP:43:GLU:HA	2.01	0.42
26:BD:10:GLY:HA2	28:BP:4:ILE:HD11	2.01	0.42
23:BB:533:G:H5'	44:BQ:23:TYR:CE2	2.54	0.42
44:BQ:26:ALA:O	44:BQ:28:SER:N	2.52	0.42
49:BR:75:VAL:HG12	49:BR:76:LYS:N	2.33	0.42
50:BT:31:VAL:C	50:BT:32:LEU:HD23	2.38	0.42
35:BV:75:GLN:HA	35:BV:75:GLN:OE1	2.19	0.42
51:BZ:70:GLU:HG3	51:BZ:71:LEU:H	1.83	0.42
1:CA:1426:G:H2'	1:CA:1427:C:C6	2.54	0.42
1:CA:415:A:O4'	1:CA:415:A:N3	2.52	0.42
1:CA:659:U:H2'	1:CA:660:C:C6	2.54	0.42
1:CA:696:A:O2'	1:CA:697:U:H5'	2.18	0.42
1:CA:782:A:H2'	1:CA:783:C:O4'	2.18	0.42
1:CA:821:G:O2'	1:CA:822:U:H5'	2.19	0.42
18:CB:74:ALA:CB	18:CB:206:ILE:HD13	2.49	0.42
18:CB:221:ARG:HA	18:CB:224:ARG:HG3	2.01	0.42
18:CB:48:MET:HB3	18:CB:199:ILE:HA	2.00	0.42
18:CB:96:LEU:H	18:CB:99:MET:HE3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:3:TYR:CE1	3:CD:10:LEU:HD21	2.54	0.42
3:CD:2:ARG:HD2	3:CD:114:ARG:HD3	2.00	0.42
1:CA:1069:C:H5''	4:CE:25:LYS:NZ	2.34	0.42
5:CF:71:ILE:O	5:CF:74:LEU:HB3	2.19	0.42
8:CI:115:VAL:HG22	8:CI:116:GLY:N	2.34	0.42
10:CK:16:SER:CA	10:CK:78:ILE:HA	2.49	0.42
11:CL:28:GLN:HB2	11:CL:80:LEU:HG	2.01	0.42
12:CM:2:ARG:HG3	12:CM:8:ILE:HG23	2.00	0.42
1:CA:473:U:OP1	13:CP:76:LYS:HE2	2.19	0.42
14:CQ:43:LEU:N	14:CQ:43:LEU:HD12	2.34	0.42
16:CS:39:ILE:HB	16:CS:66:VAL:CA	2.36	0.42
17:CT:78:LEU:O	17:CT:82:ILE:HG23	2.19	0.42
19:CU:8:ASN:O	19:CU:9:GLU:HB3	2.20	0.42
23:DB:1175:A:H2'	23:DB:1176:U:C4'	2.49	0.42
23:DB:1344:U:H5'	23:DB:1384:A:N1	2.34	0.42
23:DB:1341:G:H3'	23:DB:1397:U:O2	2.18	0.42
23:DB:163:C:O2	23:DB:163:C:O4'	2.36	0.42
23:DB:1717:A:H2'	23:DB:1718:G:O4'	2.18	0.42
23:DB:2093:G:H1'	23:DB:2198:A:C2	2.54	0.42
23:DB:248:G:N3	23:DB:2431:U:H4'	2.34	0.42
23:DB:2655:G:N2	23:DB:2664:G:H2'	2.33	0.42
23:DB:388:G:N7	23:DB:390:U:H2'	2.34	0.42
23:DB:545:U:C5	23:DB:546:U:H1'	2.54	0.42
23:DB:705:A:H61	23:DB:726:G:H1'	1.84	0.42
25:DC:109:LEU:H	25:DC:109:LEU:CD2	2.32	0.42
25:DC:157:ALA:C	25:DC:159:THR:H	2.22	0.42
26:DD:117:GLY:O	26:DD:118:PHE:C	2.57	0.42
26:DD:148:GLN:HB2	26:DD:149:ASN:H	1.64	0.42
26:DD:38:LYS:HD3	26:DD:45:TYR:CZ	2.54	0.42
48:DG:93:TYR:HA	48:DG:106:LEU:HA	2.01	0.42
48:DG:6:ALA:HB3	48:DG:68:ARG:CG	2.41	0.42
40:DH:113:SER:H	40:DH:132:PHE:HZ	1.67	0.42
24:DI:52:LEU:HD13	24:DI:81:LYS:HZ3	1.82	0.42
41:DJ:58:ASN:CA	41:DJ:127:GLY:HA2	2.49	0.42
27:DK:119:ALA:O	27:DK:120:PRO:C	2.58	0.42
23:DB:832:U:OP1	37:DL:39:LYS:HG2	2.19	0.42
37:DL:79:LEU:N	37:DL:113:ALA:CB	2.82	0.42
42:DN:17:ARG:C	42:DN:19:ALA:H	2.21	0.42
43:DO:52:SER:OG	43:DO:54:VAL:HG12	2.18	0.42
27:DK:64:ARG:NH2	28:DP:67:GLU:HG3	2.34	0.42
49:DR:15:SER:H	49:DR:18:GLN:CG	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DS:8:ARG:HA	45:DS:102:HIS:ND1	2.34	0.42
50:DT:54:GLU:CB	50:DT:88:LYS:HB2	2.44	0.42
46:DU:40:LEU:HA	46:DU:60:LYS:O	2.18	0.42
46:DU:46:LYS:HG2	46:DU:47:PRO:HD2	2.01	0.42
46:DU:34:ILE:HG23	46:DU:63:ALA:HB2	2.01	0.42
1:AA:1184:G:C2	1:AA:1185:G:C8	3.07	0.42
1:AA:1268:G:H2'	1:AA:1269:A:C8	2.53	0.42
1:AA:1329:A:O2'	1:AA:1330:U:H5'	2.20	0.42
1:AA:1385:G:H2'	1:AA:1386:G:O4'	2.19	0.42
1:AA:139:A:O2'	1:AA:140:U:H5'	2.19	0.42
1:AA:696:A:H2'	1:AA:697:U:C6	2.54	0.42
1:AA:835:U:O2'	1:AA:836:G:H5'	2.19	0.42
1:AA:967:C:H3'	1:AA:968:A:C5'	2.44	0.42
3:AD:39:GLN:HG3	3:AD:40:HIS:N	2.34	0.42
5:AF:3:HIS:CE1	5:AF:94:HIS:HA	2.54	0.42
6:AG:119:LEU:HG	6:AG:123:LEU:HD11	2.01	0.42
7:AH:86:LYS:HB3	7:AH:90:GLU:CB	2.42	0.42
1:AA:779:C:H4'	10:AK:123:PRO:HA	2.01	0.42
11:AL:18:SER:OG	11:AL:19:ASN:N	2.52	0.42
16:AS:31:ARG:HG3	16:AS:56:HIS:CE1	2.54	0.42
1:AA:1315:U:H5	16:AS:5:LYS:HZ1	1.67	0.42
32:B4:30:GLU:HA	32:B4:31:PRO:HD3	1.89	0.42
23:BB:1017:G:O2'	23:BB:1018:U:H5'	2.18	0.42
23:BB:1273:U:H4'	23:BB:1275:A:OP2	2.20	0.42
23:BB:1432:G:H2'	23:BB:1433:A:C8	2.54	0.42
23:BB:1722:A:H2'	23:BB:1723:G:C8	2.53	0.42
23:BB:2021:C:P	31:B0:8:THR:HG21	2.59	0.42
23:BB:251:A:H2'	23:BB:252:G:O4'	2.19	0.42
23:BB:2603:G:O2'	23:BB:2604:U:H5'	2.19	0.42
23:BB:2734:A:H2'	23:BB:2735:G:C5'	2.48	0.42
23:BB:2815:C:O2'	23:BB:2816:G:H5'	2.20	0.42
23:BB:2877:G:O2'	23:BB:2878:U:H5'	2.19	0.42
23:BB:345:A:H1'	23:BB:346:A:H2	1.80	0.42
23:BB:389:G:C8	23:BB:2413:G:H4'	2.54	0.42
23:BB:543:G:C2'	23:BB:544:C:H5''	2.42	0.42
23:BB:660:C:H2'	23:BB:661:A:C8	2.54	0.42
25:BC:16:VAL:CB	25:BC:203:VAL:HG11	2.46	0.42
26:BD:117:GLY:O	26:BD:118:PHE:C	2.57	0.42
48:BG:116:LEU:HD23	48:BG:120:ILE:HD13	2.01	0.42
48:BG:26:LYS:CB	48:BG:32:LEU:HG	2.49	0.42
48:BG:89:VAL:HB	48:BG:160:GLY:H	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BH:68:ARG:HH11	40:BH:72:ILE:HB	1.83	0.42
24:BI:11:GLN:NE2	24:BI:74:PRO:HG2	2.34	0.42
24:BI:126:ARG:HA	24:BI:129:GLU:OE2	2.18	0.42
41:BJ:13:ARG:HB3	41:BJ:53:TYR:HD2	1.83	0.42
41:BJ:14:ASP:O	41:BJ:53:TYR:N	2.51	0.42
27:BK:119:ALA:O	27:BK:120:PRO:O	2.37	0.42
27:BK:29:HIS:O	27:BK:30:ARG:C	2.57	0.42
27:BK:88:ASN:HD22	27:BK:89:ASN:H	1.62	0.42
42:BN:31:HIS:N	42:BN:31:HIS:ND1	2.67	0.42
43:BO:51:ALA:HB3	43:BO:78:VAL:HG13	2.01	0.42
44:BQ:50:ARG:HD2	44:BQ:50:ARG:N	2.35	0.42
46:BU:13:LEU:HD12	46:BU:13:LEU:N	2.34	0.42
35:BV:69:GLU:C	35:BV:70:ILE:HD13	2.40	0.42
52:BW:31:LEU:O	52:BW:32:ALA:HB2	2.19	0.42
52:BW:76:ARG:HH21	52:BW:76:ARG:HA	1.81	0.42
39:BX:6:LEU:N	39:BX:6:LEU:HD22	2.34	0.42
30:BY:47:ILE:CD1	30:BY:54:VAL:HG21	2.50	0.42
51:BZ:32:ASN:O	51:BZ:33:LEU:O	2.37	0.42
51:BZ:7:VAL:HG11	51:BZ:51:VAL:HG13	2.01	0.42
1:CA:1085:U:H3'	1:CA:1086:U:C6	2.53	0.42
1:CA:113:G:O4'	1:CA:354:G:H4'	2.18	0.42
1:CA:1350:A:OP2	8:CI:119:LYS:HD2	2.18	0.42
1:CA:153:C:H2'	1:CA:154:U:H6	1.84	0.42
1:CA:190:A:C4	1:CA:191:G:H1'	2.54	0.42
2:CC:155:ARG:HD2	2:CC:155:ARG:HA	1.81	0.42
2:CC:21:TRP:CH2	2:CC:31:ASN:HB3	2.55	0.42
3:CD:167:PRO:CG	3:CD:170:LEU:HD11	2.50	0.42
5:CF:21:MET:HB3	5:CF:25:TYR:CZ	2.54	0.42
9:CJ:67:ILE:HA	21:CN:94:GLY:O	2.18	0.42
10:CK:63:GLN:HG3	10:CK:98:ALA:HB1	2.00	0.42
11:CL:24:GLU:HB2	11:CL:26:CYS:HG	1.83	0.42
16:CS:28:LYS:H	16:CS:28:LYS:HD2	1.83	0.42
17:CT:64:GLY:O	17:CT:65:LEU:HB3	2.18	0.42
33:D1:38:PHE:HB2	33:D1:45:HIS:CE1	2.54	0.42
33:D1:52:LYS:CB	33:D1:52:LYS:HZ2	2.29	0.42
36:D2:22:MET:HG2	36:D2:31:LEU:HD12	2.00	0.42
22:DA:94:A:OP1	35:DV:19:ARG:HD3	2.19	0.42
23:DB:978:G:O4'	23:DB:1001:A:H2	2.02	0.42
23:DB:1028:A:N3	23:DB:2486:C:O2'	2.43	0.42
23:DB:1429:G:O2'	23:DB:1430:G:H5'	2.18	0.42
23:DB:1824:G:H2'	23:DB:1825:U:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1952:A:H5'	27:DK:42:THR:CG2	2.49	0.42
23:DB:2297:A:N6	23:DB:2319:G:H1'	2.34	0.42
23:DB:2603:G:O2'	23:DB:2604:U:H5'	2.19	0.42
23:DB:2626:C:H2'	23:DB:2627:G:H8	1.81	0.42
23:DB:327:G:O2'	23:DB:328:U:H5'	2.18	0.42
23:DB:336:C:O2'	23:DB:337:C:H5'	2.19	0.42
23:DB:481:G:C2	23:DB:507:A:C4	3.07	0.42
23:DB:808:G:H2'	23:DB:809:G:O4'	2.18	0.42
25:DC:149:LYS:HG2	25:DC:152:GLN:NE2	2.34	0.42
25:DC:151:GLY:C	25:DC:152:GLN:HE21	2.23	0.42
25:DC:68:ARG:HE	25:DC:190:THR:HG22	1.83	0.42
26:DD:159:LYS:HZ3	26:DD:159:LYS:HA	1.84	0.42
26:DD:90:PHE:O	26:DD:91:THR:C	2.57	0.42
29:DE:60:TRP:CE3	29:DE:62:GLN:HA	2.54	0.42
47:DF:21:TYR:CD2	47:DF:27:VAL:HG12	2.54	0.42
47:DF:21:TYR:HD2	47:DF:27:VAL:HG12	1.84	0.42
48:DG:96:ALA:CB	48:DG:103:ASN:HB3	2.43	0.42
48:DG:84:LYS:O	48:DG:85:LYS:O	2.37	0.42
24:DI:27:LEU:HB2	24:DI:32:VAL:HG21	2.01	0.42
24:DI:5:GLN:HB2	24:DI:30:GLN:OE1	2.19	0.42
27:DK:19:VAL:CG1	27:DK:43:ILE:HA	2.43	0.42
38:DM:21:ALA:CB	38:DM:100:LYS:HG2	2.49	0.42
38:DM:19:GLY:N	38:DM:38:ARG:HH22	2.17	0.42
38:DM:41:LEU:C	38:DM:43:ALA:H	2.22	0.42
38:DM:82:MET:HE3	38:DM:83:GLY:N	2.31	0.42
42:DN:61:ALA:C	42:DN:63:ARG:H	2.23	0.42
43:DO:49:VAL:HG23	43:DO:78:VAL:HG12	2.00	0.42
43:DO:66:GLY:C	43:DO:68:LYS:H	2.22	0.42
23:DB:560:C:H1'	44:DQ:51:GLN:HE22	1.83	0.42
49:DR:38:VAL:HB	49:DR:59:ILE:HD11	2.01	0.42
45:DS:9:HIS:H	45:DS:102:HIS:CE1	2.37	0.42
46:DU:23:LYS:HD2	46:DU:23:LYS:N	2.34	0.42
30:DY:47:ILE:CD1	30:DY:54:VAL:HG21	2.49	0.42
1:AA:1030:U:OP2	1:AA:1031:C:N4	2.52	0.42
1:AA:1347:G:N2	1:AA:1373:G:H2'	2.33	0.42
1:AA:1469:C:H3'	1:AA:1470:U:H6	1.82	0.42
1:AA:1482:G:O5'	1:AA:1482:G:H8	2.02	0.42
1:AA:223:A:O2'	1:AA:224:U:H5'	2.20	0.42
1:AA:401:C:H2'	1:AA:402:G:H8	1.84	0.42
1:AA:441:A:H61	1:AA:493:A:H62	1.67	0.42
1:AA:865:A:H5'	1:AA:1078:U:C5	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:985:C:H2'	1:AA:986:U:H6	1.84	0.42
3:AD:18:LEU:HD22	3:AD:18:LEU:N	2.34	0.42
5:AF:61:LEU:HB3	5:AF:62:MET:H	1.70	0.42
1:AA:779:C:H5''	10:AK:123:PRO:HB3	2.01	0.42
12:AM:21:ILE:CB	12:AM:24:VAL:HG22	2.46	0.42
12:AM:22:TYR:HD1	12:AM:65:GLU:HA	1.84	0.42
12:AM:90:HIS:CE1	12:AM:96:VAL:HG21	2.54	0.42
21:AN:70:HIS:O	21:AN:72:PHE:N	2.45	0.42
13:AP:67:ILE:HG12	13:AP:72:ALA:CB	2.46	0.42
19:AU:47:ALA:O	19:AU:51:ALA:N	2.52	0.42
23:BB:1050:A:N3	23:BB:1050:A:H2'	2.34	0.42
23:BB:1438:U:C4	23:BB:1552:A:N6	2.87	0.42
23:BB:1552:A:C2'	23:BB:1553:A:H5'	2.48	0.42
23:BB:1654:A:O2'	26:BD:118:PHE:HB3	2.19	0.42
23:BB:1669:A:N3	23:BB:1669:A:H2'	2.34	0.42
23:BB:1952:A:H5'	27:BK:42:THR:HG21	2.02	0.42
23:BB:2026:U:H2'	23:BB:2027:G:H8	1.83	0.42
23:BB:2180:U:O5'	23:BB:2180:U:H6	2.01	0.42
23:BB:2348:U:O2'	23:BB:2349:G:H5'	2.18	0.42
23:BB:2751:G:N3	23:BB:2751:G:H2'	2.34	0.42
23:BB:2848:G:N1	23:BB:2867:G:C2	2.87	0.42
23:BB:622:G:O2'	23:BB:623:C:H5'	2.19	0.42
23:BB:705:A:N6	23:BB:726:G:O2'	2.52	0.42
25:BC:204:LEU:HD23	25:BC:210:ALA:N	2.34	0.42
23:BB:784:G:H5''	25:BC:225:ASN:OD1	2.18	0.42
26:BD:62:LYS:HD2	26:BD:62:LYS:H	1.85	0.42
26:BD:68:PHE:CB	26:BD:73:VAL:HG23	2.47	0.42
47:BF:131:VAL:O	47:BF:133:GLU:N	2.51	0.42
47:BF:9:ASP:N	47:BF:9:ASP:OD2	2.52	0.42
48:BG:149:ALA:C	48:BG:151:ARG:H	2.23	0.42
40:BH:31:VAL:CB	40:BH:32:PRO:CD	2.88	0.42
24:BI:109:ALA:HB1	24:BI:124:MET:CG	2.49	0.42
24:BI:52:LEU:O	24:BI:54:ILE:HG13	2.19	0.42
27:BK:19:VAL:CG1	27:BK:41:ILE:HG12	2.48	0.42
27:BK:68:GLY:CA	27:BK:78:ARG:HB3	2.49	0.42
22:BA:7:G:H5''	43:BO:29:HIS:CD2	2.54	0.42
44:BQ:16:ILE:C	44:BQ:18:LYS:N	2.71	0.42
44:BQ:91:ARG:HE	44:BQ:93:ILE:CG2	2.31	0.42
46:BU:50:ALA:H	46:BU:53:GLN:NE2	2.16	0.42
46:BU:73:ASN:ND2	46:BU:74:ALA:N	2.66	0.42
35:BV:29:ILE:HD13	35:BV:31:TYR:HE2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BV:60:VAL:O	35:BV:61:LEU:HD23	2.20	0.42
35:BV:63:ILE:HB	35:BV:70:ILE:CG1	2.49	0.42
52:BW:55:ASP:O	52:BW:57:THR:N	2.50	0.42
52:BW:58:LEU:HD22	52:BW:58:LEU:N	2.33	0.42
52:BW:66:VAL:HA	52:BW:81:ILE:HG22	2.01	0.42
51:BZ:68:LEU:O	51:BZ:69:ALA:C	2.57	0.42
51:BZ:70:GLU:O	51:BZ:71:LEU:C	2.57	0.42
1:CA:1240:U:H3'	1:CA:1241:G:C5'	2.50	0.42
1:CA:1283:U:H2'	1:CA:1284:C:C6	2.54	0.42
1:CA:1409:C:H2'	1:CA:1410:A:C8	2.54	0.42
1:CA:1438:G:O2'	1:CA:1439:G:H5'	2.19	0.42
1:CA:1469:C:H3'	1:CA:1470:U:H6	1.83	0.42
1:CA:274:A:H4'	1:CA:275:G:O5'	2.19	0.42
1:CA:398:U:H2'	1:CA:399:G:C8	2.54	0.42
1:CA:517:G:N2	1:CA:533:A:OP2	2.53	0.42
1:CA:912:C:O2'	1:CA:913:A:H5'	2.19	0.42
1:CA:947:G:H2'	1:CA:948:C:H6	1.83	0.42
18:CB:150:ILE:O	18:CB:151:LYS:C	2.57	0.42
2:CC:125:ARG:HG2	2:CC:125:ARG:O	2.19	0.42
2:CC:192:TYR:N	2:CC:192:TYR:CD1	2.86	0.42
3:CD:145:ARG:O	3:CD:149:LYS:N	2.52	0.42
3:CD:178:GLU:HB3	3:CD:179:GLY:H	1.55	0.42
3:CD:191:SER:O	3:CD:192:ALA:HB2	2.18	0.42
5:CF:18:VAL:HB	5:CF:19:PRO:CD	2.49	0.42
6:CG:145:GLU:C	6:CG:147:ASN:N	2.72	0.42
6:CG:89:GLU:OE2	6:CG:89:GLU:N	2.53	0.42
7:CH:86:LYS:CD	7:CH:90:GLU:HG2	2.49	0.42
10:CK:63:GLN:HG3	10:CK:98:ALA:CB	2.49	0.42
12:CM:43:LYS:N	12:CM:43:LYS:HD2	2.34	0.42
21:CN:68:ARG:HA	21:CN:69:PRO:HD2	1.90	0.42
14:CQ:82:VAL:C	14:CQ:83:LEU:HD22	2.40	0.42
34:D3:44:ARG:N	34:D3:45:PRO:CD	2.81	0.42
23:DB:1129:A:N6	23:DB:2491:U:H5''	2.34	0.42
23:DB:1427:A:H4'	23:DB:1428:C:O4'	2.19	0.42
23:DB:1729:U:H5'	23:DB:1730:C:OP2	2.19	0.42
23:DB:1827:U:C2'	23:DB:1828:G:H5'	2.49	0.42
23:DB:1873:G:H2'	23:DB:1874:C:H6	1.85	0.42
23:DB:1977:A:H2'	23:DB:1978:A:H8	1.84	0.42
23:DB:2155:U:H2'	23:DB:2156:G:O4'	2.19	0.42
23:DB:2320:U:O2	23:DB:2320:U:H2'	2.18	0.42
23:DB:2362:C:O2'	23:DB:2363:G:H5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2508:G:H2'	23:DB:2509:G:H8	1.84	0.42
23:DB:2547:A:H4'	27:DK:29:HIS:CE1	2.54	0.42
23:DB:2700:A:O2'	23:DB:2701:U:H5'	2.19	0.42
23:DB:2688:G:H1'	23:DB:2721:A:N6	2.34	0.42
23:DB:2895:G:H2'	23:DB:2896:C:H6	1.84	0.42
23:DB:346:A:O4'	23:DB:346:A:N3	2.52	0.42
23:DB:564:C:O2'	23:DB:565:C:H5'	2.19	0.42
23:DB:581:C:H2'	23:DB:582:A:H8	1.79	0.42
23:DB:709:U:O5'	23:DB:709:U:H6	2.02	0.42
23:DB:718:A:H5'	23:DB:719:C:OP2	2.18	0.42
23:DB:903:C:H2'	23:DB:904:G:C8	2.54	0.42
23:DB:959:A:H2'	23:DB:960:A:C8	2.54	0.42
25:DC:185:ALA:C	25:DC:187:CYS:H	2.21	0.42
26:DD:173:GLN:O	26:DD:174:SER:HB3	2.18	0.42
29:DE:15:SER:C	29:DE:17:THR:H	2.23	0.42
47:DF:107:VAL:O	47:DF:110:ILE:HG22	2.19	0.42
41:DJ:29:ALA:O	41:DJ:32:LEU:HB2	2.19	0.42
37:DL:143:GLU:CG	37:DL:144:GLU:N	2.77	0.42
38:DM:26:VAL:HB	38:DM:104:GLU:OE2	2.19	0.42
38:DM:126:ILE:HG22	38:DM:127:LYS:N	2.34	0.42
38:DM:19:GLY:N	38:DM:38:ARG:NH1	2.65	0.42
42:DN:71:ARG:HG2	42:DN:71:ARG:HH21	1.84	0.42
43:DO:7:ARG:HG3	43:DO:96:GLY:C	2.39	0.42
44:DQ:86:SER:HB3	49:DR:50:GLY:O	2.19	0.42
44:DQ:90:ASP:H	49:DR:39:LEU:HD11	1.83	0.42
45:DS:45:VAL:HG23	45:DS:46:LEU:HD23	2.00	0.42
50:DT:51:PHE:HB3	50:DT:53:VAL:HG23	2.01	0.42
35:DV:38:LEU:HD11	35:DV:40:ILE:HG23	2.01	0.42
52:DW:77:LYS:HA	52:DW:77:LYS:HD3	1.81	0.42
1:AA:1085:U:O4'	1:AA:1094:G:N1	2.52	0.42
1:AA:1342:C:O2'	8:AI:125:GLN:CB	2.68	0.42
1:AA:1453:G:N3	1:AA:1453:G:H3'	2.34	0.42
1:AA:1479:C:O2'	1:AA:1480:A:H5'	2.19	0.42
1:AA:490:C:H2'	1:AA:491:G:O4'	2.19	0.42
1:AA:633:G:H2'	1:AA:634:C:C6	2.55	0.42
1:AA:769:G:H4'	1:AA:1513:A:H4'	2.01	0.42
1:AA:904:U:H2'	1:AA:905:U:C6	2.54	0.42
1:AA:904:U:H6	1:AA:904:U:O5'	2.02	0.42
3:AD:145:ARG:O	3:AD:149:LYS:N	2.52	0.42
4:AE:93:VAL:HA	4:AE:126:ALA:CB	2.50	0.42
6:AG:87:PRO:CB	6:AG:144:ALA:HA	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:74:VAL:HB	6:AG:85:GLN:HB3	2.01	0.42
7:AH:100:ILE:HD11	7:AH:102:VAL:HG22	2.01	0.42
8:AI:71:ILE:N	8:AI:71:ILE:HD12	2.28	0.42
10:AK:34:THR:HB	10:AK:40:ALA:HA	2.01	0.42
11:AL:32:VAL:O	11:AL:33:CYS:HB3	2.20	0.42
21:AN:52:ARG:HB3	21:AN:53:ASP:H	1.56	0.42
20:AO:81:LEU:HD23	20:AO:81:LEU:C	2.40	0.42
20:AO:8:THR:O	20:AO:11:ILE:HG22	2.19	0.42
14:AQ:10:ARG:O	14:AQ:22:VAL:HG13	2.20	0.42
15:AR:61:ALA:HB3	15:AR:67:LEU:HD12	2.02	0.42
17:AT:44:ALA:O	17:AT:48:LYS:HB3	2.19	0.42
19:AU:20:ARG:H	19:AU:20:ARG:CD	2.32	0.42
33:B1:4:ILE:HB	33:B1:27:ARG:NH1	2.35	0.42
22:BA:67:G:H2'	22:BA:68:C:H6	1.84	0.42
23:BB:1799:G:OP1	25:BC:257:ARG:HG3	2.19	0.42
23:BB:1909:C:O2'	23:BB:1910:G:H5'	2.20	0.42
23:BB:2425:A:H4'	23:BB:2426:A:H5''	2.02	0.42
23:BB:2552:U:H3'	23:BB:2554:U:OP2	2.19	0.42
23:BB:255:A:H2'	23:BB:256:A:O4'	2.19	0.42
23:BB:271:G:N2	23:BB:367:G:H1'	2.35	0.42
23:BB:478:A:H5''	23:BB:479:A:OP2	2.19	0.42
23:BB:819:A:OP2	23:BB:1187:G:N2	2.53	0.42
23:BB:83:A:H61	23:BB:101:A:C5'	2.31	0.42
23:BB:8:C:H5''	41:BJ:53:TYR:OH	2.19	0.42
23:BB:97:C:H2'	23:BB:98:G:O4'	2.19	0.42
25:BC:183:VAL:HG22	25:BC:184:GLU:H	1.83	0.42
26:BD:38:LYS:HD3	26:BD:45:TYR:CZ	2.55	0.42
29:BE:109:LEU:O	29:BE:112:LEU:HB2	2.19	0.42
29:BE:69:ARG:O	29:BE:70:SER:CB	2.66	0.42
47:BF:3:LEU:O	47:BF:6:TYR:HB3	2.19	0.42
48:BG:74:MET:HG3	48:BG:74:MET:H	1.59	0.42
24:BI:103:ALA:O	24:BI:107:GLU:HG3	2.20	0.42
37:BL:79:LEU:HB2	37:BL:113:ALA:N	2.32	0.42
38:BM:69:PRO:HA	38:BM:93:VAL:O	2.20	0.42
41:BJ:40:HIS:O	44:BQ:66:ALA:HB1	2.19	0.42
44:BQ:63:ARG:HH22	44:BQ:96:ASP:N	2.18	0.42
50:BT:54:GLU:O	50:BT:55:VAL:HB	2.19	0.42
1:CA:1026:G:H22	1:CA:1035:A:H2	1.65	0.42
1:CA:1387:G:H2'	1:CA:1388:C:H6	1.84	0.42
1:CA:28:A:N3	1:CA:296:U:H4'	2.34	0.42
1:CA:34:C:H2'	1:CA:35:G:H8	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:667:G:H2'	1:CA:668:G:H8	1.84	0.42
1:CA:79:G:O2'	1:CA:80:A:H5'	2.19	0.42
1:CA:824:G:C6	1:CA:877:G:C6	3.08	0.42
1:CA:835:U:O2'	1:CA:836:G:H5'	2.20	0.42
6:CG:31:VAL:HG22	6:CG:32:ASP:OD2	2.20	0.42
6:CG:90:VAL:HG23	6:CG:91:ARG:O	2.19	0.42
7:CH:100:ILE:HD11	7:CH:102:VAL:HG22	2.02	0.42
8:CI:29:ILE:HA	8:CI:64:ILE:O	2.19	0.42
9:CJ:49:PHE:O	9:CJ:64:GLN:HA	2.19	0.42
9:CJ:5:ARG:N	9:CJ:76:ILE:O	2.52	0.42
10:CK:85:VAL:HG12	10:CK:86:LYS:N	2.34	0.42
12:CM:2:ARG:HD3	12:CM:2:ARG:H	1.84	0.42
21:CN:50:LEU:N	21:CN:51:PRO:CD	2.75	0.42
13:CP:39:PHE:CG	13:CP:40:ASN:N	2.87	0.42
36:D2:30:VAL:O	36:D2:33:ARG:HB2	2.20	0.42
23:DB:1128:G:C6	23:DB:2518:A:N6	2.87	0.42
23:DB:1214:A:H2'	23:DB:1215:G:O4'	2.18	0.42
23:DB:1524:G:H2'	23:DB:1525:A:O4'	2.19	0.42
23:DB:1700:A:H2'	23:DB:1701:A:C5'	2.49	0.42
23:DB:1877:A:H2'	23:DB:1878:G:C8	2.55	0.42
23:DB:1917:U:H2'	23:DB:1918:A:C5'	2.46	0.42
23:DB:1853:A:N1	23:DB:2087:G:H1'	2.34	0.42
23:DB:2420:C:OP2	34:D3:32:LEU:N	2.52	0.42
23:DB:265:A:O2'	23:DB:266:G:H4'	2.18	0.42
23:DB:2795:C:H2'	23:DB:2796:U:C4'	2.50	0.42
23:DB:730:A:H3'	56:DB:3609:HOH:O	2.20	0.42
25:DC:75:ALA:HA	25:DC:96:LYS:NZ	2.35	0.42
26:DD:114:LYS:HD2	26:DD:116:LYS:HE3	2.01	0.42
26:DD:2:ILE:O	26:DD:2:ILE:HD12	2.20	0.42
47:DF:121:PHE:CA	47:DF:127:TYR:HA	2.48	0.42
47:DF:28:PRO:HB2	47:DF:168:LEU:CD2	2.40	0.42
48:DG:148:ARG:HH11	48:DG:152:ARG:NE	2.17	0.42
48:DG:16:VAL:CG1	48:DG:49:LEU:HD13	2.49	0.42
40:DH:95:GLY:O	40:DH:99:ILE:HG12	2.19	0.42
49:DR:3:ALA:HB2	49:DR:101:ILE:HD12	2.00	0.42
50:DT:39:THR:HG22	50:DT:42:GLU:CG	2.48	0.42
52:DW:67:LYS:O	52:DW:68:PHE:HB2	2.19	0.42
1:AA:1363:A:H2'	1:AA:1363:A:N3	2.35	0.42
1:AA:1381:U:H2'	1:AA:1382:C:C6	2.55	0.42
1:AA:274:A:H4'	1:AA:275:G:O5'	2.19	0.42
1:AA:321:A:O2'	1:AA:322:C:H5'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:343:U:O2'	1:AA:344:A:H8	2.03	0.42
1:AA:382:A:H2'	1:AA:383:A:H8	1.82	0.42
1:AA:492:C:H2'	1:AA:493:A:N3	2.34	0.42
1:AA:509:A:C6	1:AA:510:A:N1	2.88	0.42
1:AA:532:A:H62	2:AC:191:THR:CB	2.32	0.42
1:AA:572:A:N3	1:AA:917:G:H1'	2.33	0.42
1:AA:725:G:H2'	1:AA:726:C:C6	2.55	0.42
1:AA:740:U:H4'	20:AO:42:HIS:CD2	2.54	0.42
1:AA:746:A:H2'	1:AA:747:A:C8	2.54	0.42
18:AB:112:ARG:NH2	18:AB:116:LEU:HD11	2.31	0.42
18:AB:69:VAL:HG12	18:AB:70:GLY:N	2.34	0.42
2:AC:52:SER:C	2:AC:113:LYS:HG2	2.39	0.42
3:AD:53:GLN:HG2	3:AD:198:LEU:HD22	2.01	0.42
3:AD:61:ARG:HG3	3:AD:71:PHE:CD2	2.54	0.42
4:AE:14:LEU:C	4:AE:14:LEU:HD13	2.40	0.42
4:AE:20:VAL:O	4:AE:31:SER:HB2	2.19	0.42
4:AE:77:ASN:HD22	4:AE:77:ASN:HA	1.64	0.42
6:AG:53:SER:O	6:AG:55:LYS:N	2.52	0.42
11:AL:27:PRO:O	11:AL:28:GLN:NE2	2.52	0.42
11:AL:85:ARG:HB2	11:AL:93:ARG:HA	2.01	0.42
15:AR:34:GLU:HB2	19:AU:18:PHE:CZ	2.55	0.42
16:AS:35:ARG:CG	16:AS:36:ARG:HH21	2.32	0.42
33:B1:35:LEU:O	33:B1:36:LYS:HB3	2.20	0.42
36:B2:30:VAL:O	36:B2:33:ARG:HB2	2.19	0.42
22:BA:112:G:H2'	22:BA:113:C:H6	1.85	0.42
22:BA:3:C:H2'	22:BA:4:C:C6	2.54	0.42
22:BA:52:A:O2'	22:BA:53:A:H5'	2.20	0.42
23:BB:1198:U:H2'	23:BB:1199:U:H6	1.83	0.42
23:BB:1407:G:O2'	23:BB:1408:G:H5'	2.20	0.42
23:BB:1649:G:O2'	23:BB:1650:A:H5'	2.20	0.42
23:BB:1727:C:H2'	23:BB:1728:C:O4'	2.19	0.42
23:BB:1930:G:H2'	23:BB:1968:G:H1	1.85	0.42
23:BB:2014:A:H2'	23:BB:2015:A:C8	2.53	0.42
23:BB:216:A:H2'	23:BB:217:A:C8	2.52	0.42
23:BB:230:G:H2'	23:BB:231:A:C8	2.55	0.42
23:BB:2358:A:H2'	23:BB:2359:C:O4'	2.20	0.42
23:BB:2688:G:H1'	23:BB:2721:A:N6	2.34	0.42
23:BB:2772:C:H2'	23:BB:2773:C:C6	2.55	0.42
23:BB:381:G:O2'	23:BB:382:A:H5'	2.20	0.42
23:BB:672:C:H2'	23:BB:673:C:H6	1.84	0.42
23:BB:707:G:O2'	23:BB:708:G:H5'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:705:A:H61	23:BB:726:G:H1'	1.81	0.42
23:BB:740:C:O2'	23:BB:741:U:H5'	2.20	0.42
23:BB:805:G:N2	23:BB:829:A:OP1	2.52	0.42
25:BC:63:ILE:HD13	25:BC:63:ILE:HA	1.82	0.42
48:BG:157:LYS:HG2	48:BG:159:LYS:CG	2.49	0.42
48:BG:169:ARG:HB3	48:BG:171:LYS:NZ	2.34	0.42
48:BG:90:GLY:HA3	48:BG:93:TYR:CE1	2.55	0.42
24:BI:63:ASP:C	24:BI:65:SER:H	2.22	0.42
27:BK:99:ILE:HG12	27:BK:115:ILE:HG13	2.01	0.42
27:BK:88:ASN:ND2	27:BK:89:ASN:N	2.62	0.42
42:BN:35:LYS:HG2	42:BN:112:TYR:CE1	2.54	0.42
43:BO:64:TYR:CB	43:BO:67:ASN:HB2	2.37	0.42
44:BQ:101:ASP:HB2	49:BR:2:TYR:OH	2.20	0.42
49:BR:58:VAL:C	49:BR:59:ILE:HG13	2.40	0.42
1:CA:1320:C:N4	16:CS:35:ARG:HG3	2.34	0.42
1:CA:16:A:C2'	1:CA:17:U:H5'	2.50	0.42
1:CA:193:C:H2'	1:CA:194:C:C5	2.54	0.42
1:CA:411:A:N6	1:CA:413:G:H21	2.17	0.42
1:CA:696:A:H2'	1:CA:697:U:H6	1.84	0.42
1:CA:707:U:H4'	10:CK:21:HIS:CG	2.54	0.42
1:CA:776:G:N2	1:CA:803:G:N7	2.67	0.42
1:CA:916:U:H2'	1:CA:917:G:H8	1.84	0.42
18:CB:137:THR:O	18:CB:140:LEU:HB2	2.19	0.42
18:CB:60:ALA:CB	18:CB:224:ARG:HG2	2.50	0.42
18:CB:58:LYS:C	18:CB:60:ALA:H	2.22	0.42
9:CJ:48:ARG:HA	9:CJ:66:GLU:HA	2.01	0.42
9:CJ:80:THR:C	9:CJ:82:LYS:H	2.22	0.42
9:CJ:9:ARG:HH11	9:CJ:9:ARG:HG2	1.84	0.42
12:CM:2:ARG:HB2	12:CM:2:ARG:HE	1.51	0.42
16:CS:47:THR:HA	16:CS:60:PHE:CE2	2.55	0.42
19:CU:36:PHE:O	19:CU:39:LYS:HD2	2.19	0.42
23:DB:1099:G:N7	24:DI:3:LYS:CD	2.83	0.42
23:DB:1344:U:H5'	23:DB:1384:A:C6	2.54	0.42
23:DB:1417:C:C4'	23:DB:1587:G:H21	2.32	0.42
23:DB:1745:A:H2'	23:DB:1746:A:C8	2.55	0.42
23:DB:1767:G:O2'	23:DB:1768:C:H5'	2.20	0.42
23:DB:1940:U:O2	23:DB:1940:U:H5''	2.20	0.42
23:DB:2305:U:H6	47:DF:152:ASP:OD2	2.02	0.42
23:DB:2346:A:O4'	23:DB:2383:G:O4'	2.36	0.42
23:DB:2648:G:H2'	23:DB:2649:C:H6	1.84	0.42
23:DB:2650:U:O2'	23:DB:2651:C:H5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:28:A:O2'	23:DB:583:G:H5'	2.20	0.42
23:DB:714:U:C2	23:DB:716:A:H3'	2.54	0.42
23:DB:950:G:H2'	23:DB:951:C:H6	1.79	0.42
25:DC:109:LEU:HD23	25:DC:109:LEU:N	2.35	0.42
26:DD:109:VAL:HG11	26:DD:193:VAL:HB	2.01	0.42
47:DF:113:PHE:HZ	47:DF:175:PRO:HB2	1.83	0.42
47:DF:37:MET:O	47:DF:37:MET:HG2	2.19	0.42
48:DG:116:LEU:HD23	48:DG:120:ILE:HD13	2.02	0.42
48:DG:83:THR:HA	48:DG:84:LYS:HZ1	1.84	0.42
48:DG:93:TYR:HD2	48:DG:93:TYR:N	2.18	0.42
40:DH:143:ILE:HG22	40:DH:144:VAL:N	2.34	0.42
40:DH:37:VAL:HG13	40:DH:37:VAL:O	2.19	0.42
40:DH:42:LYS:HD3	40:DH:42:LYS:C	2.39	0.42
41:DJ:25:LEU:HB2	41:DJ:62:VAL:HG22	2.00	0.42
38:DM:29:GLY:CA	38:DM:106:ASP:HB2	2.49	0.42
43:DO:7:ARG:O	43:DO:10:ARG:HB2	2.19	0.42
44:DQ:26:ALA:C	44:DQ:28:SER:N	2.73	0.42
49:DR:27:ILE:HA	49:DR:27:ILE:HD13	1.85	0.42
50:DT:15:HIS:H	50:DT:32:LEU:CA	2.21	0.42
23:DB:2365:G:H4'	52:DW:59:PHE:CD1	2.54	0.42
52:DW:8:SER:O	52:DW:9:THR:HB	2.18	0.42
39:DX:7:ARG:HA	39:DX:7:ARG:NE	2.35	0.42
1:AA:1048:G:N2	1:AA:1214:C:C5	2.75	0.42
1:AA:1286:U:H2'	1:AA:1286:U:H6	1.68	0.42
1:AA:140:U:H2'	1:AA:141:G:C8	2.55	0.42
18:AB:166:ASP:CG	18:AB:190:SER:HA	2.40	0.42
2:AC:130:ARG:O	2:AC:133:MET:HB2	2.19	0.42
2:AC:167:TYR:CD2	2:AC:168:ARG:N	2.87	0.42
2:AC:40:GLN:O	2:AC:43:THR:HB	2.20	0.42
3:AD:181:PHE:O	3:AD:182:LYS:C	2.57	0.42
3:AD:61:ARG:NH1	3:AD:68:GLU:HB2	2.34	0.42
3:AD:78:ALA:C	3:AD:85:THR:HG23	2.39	0.42
5:AF:29:ILE:HG21	5:AF:64:VAL:CG1	2.47	0.42
7:AH:74:ILE:HG12	7:AH:74:ILE:O	2.20	0.42
10:AK:115:ILE:HG21	19:AU:23:GLU:OE1	2.20	0.42
1:AA:529:G:O6	11:AL:45:ASN:HA	2.20	0.42
21:AN:33:VAL:HA	21:AN:40:ARG:HB2	2.01	0.42
21:AN:58:ARG:C	21:AN:59:GLN:HE21	2.22	0.42
13:AP:33:ILE:HG13	13:AP:33:ILE:H	1.69	0.42
13:AP:36:VAL:HG13	13:AP:36:VAL:O	2.19	0.42
1:AA:720:C:H5"	15:AR:40:PRO:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:B1:8:ILE:HD12	33:B1:51:ALA:HA	2.01	0.42
23:BB:651:G:O5'	34:B3:17:GLY:HA3	2.19	0.42
23:BB:1051:G:H2'	23:BB:1052:C:H6	1.82	0.42
23:BB:1125:G:C6	23:BB:1126:A:N6	2.88	0.42
23:BB:1153:C:H2'	23:BB:1154:G:C8	2.54	0.42
23:BB:1241:A:H3'	23:BB:1242:U:H6	1.84	0.42
23:BB:1319:C:C2'	23:BB:1320:C:H5'	2.49	0.42
23:BB:1400:U:O2'	23:BB:1401:G:H5'	2.19	0.42
23:BB:1684:G:H2'	23:BB:1685:C:H6	1.85	0.42
23:BB:1750:G:O2'	23:BB:1751:U:H5'	2.20	0.42
23:BB:175:G:O2'	23:BB:176:A:H5'	2.20	0.42
23:BB:1977:A:H2'	23:BB:1978:A:H8	1.85	0.42
23:BB:2386:A:H2'	23:BB:2387:U:C6	2.55	0.42
23:BB:957:C:N4	23:BB:2459:A:C8	2.87	0.42
23:BB:2630:G:H2'	23:BB:2631:G:C8	2.49	0.42
23:BB:2650:U:O2'	23:BB:2651:C:H5'	2.20	0.42
23:BB:460:A:H2'	23:BB:461:C:O4'	2.20	0.42
23:BB:488:G:N2	23:BB:493:G:O6	2.53	0.42
23:BB:562:U:C4	23:BB:2036:C:O4'	2.72	0.42
23:BB:986:C:O2'	23:BB:987:C:H5'	2.20	0.42
25:BC:202:ARG:HH11	25:BC:213:ARG:NE	2.18	0.42
25:BC:61:TYR:HA	25:BC:85:ASN:ND2	2.34	0.42
26:BD:121:THR:C	26:BD:123:LYS:N	2.71	0.42
26:BD:130:GLN:HB3	26:BD:130:GLN:HE21	1.57	0.42
26:BD:36:GLN:HE21	26:BD:36:GLN:HB2	1.62	0.42
26:BD:46:ARG:HH11	26:BD:46:ARG:HG3	1.84	0.42
29:BE:173:THR:HA	29:BE:199:MET:HE1	2.01	0.42
47:BF:78:ILE:C	47:BF:79:ARG:HG3	2.40	0.42
48:BG:16:VAL:CG1	48:BG:49:LEU:HD13	2.50	0.42
48:BG:8:VAL:HB	48:BG:49:LEU:O	2.20	0.42
37:BL:85:VAL:O	37:BL:85:VAL:HG22	2.20	0.42
38:BM:110:GLU:HG2	38:BM:114:ARG:HH12	1.84	0.42
38:BM:71:LYS:CG	38:BM:93:VAL:HG12	2.49	0.42
42:BN:55:ALA:HB1	42:BN:84:GLY:HA2	2.01	0.42
43:BO:7:ARG:O	43:BO:10:ARG:HB2	2.20	0.42
28:BP:57:ALA:O	28:BP:58:PHE:HD2	2.02	0.42
23:BB:560:C:H1'	44:BQ:51:GLN:HE22	1.84	0.42
46:BU:21:ARG:HD2	46:BU:72:PHE:CD2	2.55	0.42
52:BW:37:VAL:C	52:BW:38:ARG:HG2	2.40	0.42
1:CA:1411:C:H2'	1:CA:1412:C:H6	1.85	0.42
1:CA:1525:G:O2'	1:CA:1526:G:H5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:229:U:H2'	1:CA:230:G:H8	1.83	0.42
1:CA:126:G:C4'	1:CA:634:C:H1'	2.49	0.42
1:CA:91:U:C2	1:CA:92:U:C5	3.08	0.42
1:CA:929:G:H2'	1:CA:930:C:O4'	2.20	0.42
18:CB:128:LEU:HD23	18:CB:133:ALA:HA	2.02	0.42
3:CD:24:VAL:HG12	3:CD:160:LEU:HB3	2.01	0.42
3:CD:4:LEU:N	3:CD:4:LEU:HD12	2.34	0.42
3:CD:53:GLN:HG2	3:CD:198:LEU:HD22	2.02	0.42
4:CE:77:ASN:HB3	4:CE:78:GLY:H	1.70	0.42
5:CF:51:ILE:HD11	5:CF:86:ARG:HG3	2.01	0.42
6:CG:72:VAL:HG12	6:CG:89:GLU:HA	2.01	0.42
7:CH:68:LYS:HD2	7:CH:69:ALA:H	1.84	0.42
12:CM:13:HIS:ND1	12:CM:41:ASP:HA	2.35	0.42
13:CP:67:ILE:HD11	13:CP:71:VAL:CG2	2.46	0.42
13:CP:72:ALA:HA	13:CP:75:ILE:CD1	2.49	0.42
33:D1:8:ILE:HB	33:D1:24:LYS:HB2	2.01	0.42
22:DA:8:C:O2'	43:DO:40:ILE:HD13	2.19	0.42
23:DB:814:C:H1'	23:DB:1225:G:N2	2.34	0.42
23:DB:1353:A:O2'	23:DB:1354:A:H5'	2.18	0.42
23:DB:1923:U:H2'	23:DB:1924:C:C6	2.54	0.42
23:DB:2011:U:H2'	23:DB:2012:G:O4'	2.20	0.42
23:DB:2060:A:H1'	56:DB:3297:HOH:O	2.19	0.42
23:DB:250:G:H2'	23:DB:251:A:C8	2.55	0.42
23:DB:348:A:H2'	23:DB:349:U:O4'	2.19	0.42
23:DB:478:A:H5''	23:DB:479:A:OP2	2.20	0.42
23:DB:513:A:O5'	23:DB:513:A:H8	2.01	0.42
25:DC:14:HIS:O	25:DC:16:VAL:HG23	2.20	0.42
25:DC:244:VAL:HG12	25:DC:250:GLN:HA	2.01	0.42
25:DC:57:HIS:ND1	25:DC:58:LYS:N	2.67	0.42
26:DD:150:GLN:O	26:DD:153:GLY:N	2.52	0.42
29:DE:152:GLU:O	29:DE:153:LEU:HB3	2.20	0.42
48:DG:97:VAL:CB	48:DG:124:CYS:HB2	2.49	0.42
48:DG:151:ARG:O	48:DG:152:ARG:HB3	2.20	0.42
48:DG:26:LYS:HA	48:DG:32:LEU:CA	2.50	0.42
48:DG:88:LEU:HD13	48:DG:93:TYR:HB2	2.01	0.42
40:DH:90:LEU:HB2	40:DH:123:ARG:HA	2.01	0.42
27:DK:60:ALA:HA	27:DK:87:LEU:CD2	2.50	0.42
27:DK:8:LEU:HD12	27:DK:8:LEU:N	2.34	0.42
37:DL:93:ASN:HD22	37:DL:94:THR:H	1.62	0.42
37:DL:95:LEU:HB2	37:DL:101:ILE:HG12	2.01	0.42
42:DN:29:VAL:O	42:DN:78:LYS:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DP:109:ILE:HG13	28:DP:109:ILE:O	2.20	0.42
44:DQ:81:GLY:HA3	44:DQ:112:ALA:HB1	2.01	0.42
35:DV:11:GLU:HB2	35:DV:12:GLN:H	1.58	0.42
51:DZ:49:LEU:O	51:DZ:51:VAL:HG23	2.19	0.42
1:AA:1198:G:H5''	56:AA:1865:HOH:O	2.19	0.42
1:AA:1047:G:H21	1:AA:1215:G:C4'	2.33	0.42
1:AA:1237:C:H5''	1:AA:1238:A:H5'	2.01	0.42
1:AA:1339:A:H2'	1:AA:1340:A:H5'	2.02	0.42
1:AA:1355:G:H2'	1:AA:1356:G:H8	1.85	0.42
1:AA:1392:G:O2'	1:AA:1393:U:H5'	2.20	0.42
1:AA:1495:U:H2'	1:AA:1496:C:C6	2.55	0.42
1:AA:1520:C:H2'	1:AA:1521:C:C6	2.55	0.42
1:AA:1525:G:O2'	1:AA:1526:G:H5'	2.19	0.42
1:AA:17:U:O4'	1:AA:1080:A:H1'	2.19	0.42
1:AA:399:G:H2'	1:AA:400:C:H6	1.84	0.42
1:AA:989:U:O2'	1:AA:990:C:H5'	2.19	0.42
18:AB:169:HIS:HA	18:AB:172:ILE:CD1	2.48	0.42
2:AC:102:ILE:HG22	2:AC:103:ALA:N	2.33	0.42
2:AC:146:LYS:HE3	2:AC:202:PHE:CE2	2.54	0.42
4:AE:45:VAL:H	4:AE:71:ILE:HG22	1.83	0.42
9:AJ:14:ASP:HA	9:AJ:70:HIS:HD2	1.85	0.42
21:AN:30:ILE:HD12	21:AN:44:VAL:HG13	2.01	0.42
22:BA:19:C:H2'	22:BA:20:G:H8	1.85	0.42
23:BB:1063:G:O2'	24:BI:88:GLY:HA3	2.19	0.42
23:BB:1291:C:O2'	23:BB:1292:G:H5'	2.20	0.42
23:BB:1399:C:H2'	23:BB:1400:U:C6	2.54	0.42
23:BB:1654:A:O2'	26:BD:118:PHE:CB	2.68	0.42
23:BB:1799:G:N2	23:BB:1818:U:O2'	2.53	0.42
23:BB:2233:U:H2'	23:BB:2234:G:C8	2.55	0.42
23:BB:244:A:OP2	34:B3:7:ARG:NH2	2.45	0.42
23:BB:2527:C:H2'	23:BB:2528:U:H6	1.85	0.42
23:BB:2636:C:H2'	23:BB:2637:U:H6	1.85	0.42
23:BB:2808:G:O2'	23:BB:2809:A:H8	2.02	0.42
23:BB:381:G:H5''	51:BZ:16:ASN:ND2	2.35	0.42
23:BB:557:C:H2'	23:BB:558:U:H6	1.82	0.42
23:BB:63:A:O5'	23:BB:63:A:H8	2.01	0.42
23:BB:754:U:H2'	23:BB:755:U:C6	2.54	0.42
23:BB:974:G:H2'	23:BB:974:G:N3	2.35	0.42
26:BD:5:VAL:HG11	26:BD:80:TRP:CE3	2.55	0.42
47:BF:110:ILE:HA	47:BF:111:ARG:NH2	2.35	0.42
47:BF:11:VAL:HG12	47:BF:12:VAL:N	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BI:78:LEU:HD13	24:BI:108:ILE:HG23	2.01	0.42
27:BK:3:GLN:CG	27:BK:4:GLU:N	2.82	0.42
37:BL:143:GLU:CG	37:BL:144:GLU:N	2.79	0.42
38:BM:73:ILE:HG21	38:BM:91:TYR:OH	2.19	0.42
23:BB:2840:C:H5''	42:BN:53:THR:HG21	2.02	0.42
28:BP:59:THR:OG1	28:BP:72:VAL:HG12	2.19	0.42
28:BP:79:VAL:HG23	28:BP:80:VAL:N	2.35	0.42
49:BR:64:VAL:O	49:BR:65:ALA:HB3	2.19	0.42
23:BB:973:A:H5'	49:BR:81:LYS:HE3	2.00	0.42
45:BS:16:LYS:O	45:BS:19:LEU:HB3	2.19	0.42
45:BS:29:VAL:O	45:BS:32:ALA:HB3	2.19	0.42
50:BT:51:PHE:C	50:BT:53:VAL:H	2.23	0.42
46:BU:12:VAL:HG21	46:BU:38:ILE:HG12	2.02	0.42
46:BU:28:LEU:HG	46:BU:34:ILE:HD11	2.02	0.42
52:BW:16:GLU:O	52:BW:17:ALA:C	2.58	0.42
1:CA:1226:C:H4'	1:CA:1227:A:OP1	2.20	0.42
1:CA:1248:A:H2'	1:CA:1249:C:C6	2.55	0.42
1:CA:134:G:H2'	1:CA:135:C:O4'	2.18	0.42
1:CA:160:A:H2'	1:CA:161:A:C8	2.55	0.42
1:CA:312:C:O2'	1:CA:313:A:H5'	2.19	0.42
1:CA:465:A:H2'	1:CA:467:U:OP1	2.20	0.42
1:CA:72:A:N6	1:CA:98:A:H2	2.18	0.42
1:CA:981:U:C5'	21:CN:60:ARG:HD2	2.50	0.42
18:CB:21:TYR:O	18:CB:22:TRP:O	2.37	0.42
18:CB:47:PRO:O	18:CB:51:GLU:HB2	2.20	0.42
2:CC:172:VAL:HG12	2:CC:174:LEU:HD12	2.02	0.42
2:CC:4:VAL:HG22	2:CC:5:HIS:N	2.34	0.42
1:CA:409:U:OP1	3:CD:21:LYS:HG3	2.20	0.42
4:CE:14:LEU:HD13	4:CE:14:LEU:C	2.40	0.42
6:CG:134:VAL:HG12	6:CG:137:ARG:NH2	2.35	0.42
6:CG:87:PRO:HG3	6:CG:148:LYS:CA	2.37	0.42
7:CH:17:GLN:HE21	7:CH:62:LEU:CD2	2.33	0.42
4:CE:158:LYS:HZ3	7:CH:63:LYS:HD3	1.83	0.42
8:CI:56:MET:O	8:CI:57:VAL:HB	2.20	0.42
10:CK:27:ASN:O	10:CK:56:LYS:HE2	2.20	0.42
21:CN:25:GLU:CD	21:CN:25:GLU:H	2.21	0.42
21:CN:86:ALA:HA	21:CN:89:ARG:NH2	2.34	0.42
13:CP:3:THR:HG22	13:CP:66:THR:HB	2.02	0.42
36:D2:10:LEU:O	36:D2:14:ARG:HG2	2.19	0.42
22:DA:63:C:H2'	22:DA:64:G:H8	1.84	0.42
23:DB:1399:C:H2'	23:DB:1400:U:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1567:G:H1'	23:DB:1568:G:C5	2.55	0.42
23:DB:1930:G:C2'	23:DB:1931:U:OP2	2.67	0.42
23:DB:2219:U:O2'	23:DB:2220:U:H5'	2.20	0.42
23:DB:2232:C:P	51:DZ:27:ARG:HH22	2.43	0.42
23:DB:2271:G:H2'	23:DB:2272:U:C5	2.54	0.42
23:DB:2387:U:H1'	52:DW:38:ARG:NH2	2.34	0.42
23:DB:367:G:N2	23:DB:368:A:H1'	2.34	0.42
23:DB:527:C:C4	23:DB:2779:U:H2'	2.54	0.42
23:DB:527:C:N4	23:DB:2779:U:OP2	2.52	0.42
23:DB:850:U:H4'	30:DY:22:THR:HG22	2.02	0.42
25:DC:209:ALA:O	25:DC:213:ARG:HB2	2.19	0.42
29:DE:146:VAL:HG12	29:DE:147:LEU:N	2.35	0.42
48:DG:26:LYS:CB	48:DG:32:LEU:HG	2.49	0.42
48:DG:29:ASN:HD22	48:DG:30:GLY:N	2.17	0.42
41:DJ:100:VAL:O	41:DJ:104:ALA:HB2	2.19	0.42
41:DJ:81:ILE:HG23	41:DJ:82:GLY:N	2.25	0.42
38:DM:65:ILE:HG23	38:DM:103:TYR:CD2	2.55	0.42
38:DM:29:GLY:HA2	38:DM:106:ASP:HB2	2.00	0.42
42:DN:33:ILE:HD12	42:DN:34:ILE:N	2.35	0.42
28:DP:48:ALA:HB3	28:DP:59:THR:OG1	2.20	0.42
28:DP:95:LYS:HG3	28:DP:97:TYR:CE1	2.54	0.42
44:DQ:16:ILE:C	44:DQ:18:LYS:N	2.73	0.42
44:DQ:26:ALA:O	44:DQ:28:SER:N	2.52	0.42
44:DQ:40:LYS:O	44:DQ:43:GLN:HB2	2.19	0.42
23:DB:992:C:H4'	49:DR:74:ILE:HD13	2.02	0.42
49:DR:79:ARG:O	49:DR:80:ARG:HB2	2.20	0.42
45:DS:4:ILE:CG2	45:DS:106:VAL:HG22	2.49	0.42
1:AA:1127:G:H5'	1:AA:1280:A:HO2'	1.85	0.42
1:AA:1212:U:H5'	1:AA:1213:A:O4'	2.20	0.42
1:AA:982:U:H3	1:AA:1223:C:H42	1.68	0.42
1:AA:1267:C:H2'	1:AA:1268:G:C8	2.55	0.42
1:AA:1360:A:N6	1:AA:1361:G:N3	2.68	0.42
1:AA:143:A:O3'	1:AA:144:G:H8	2.03	0.42
1:AA:144:G:H2'	1:AA:145:G:O4'	2.20	0.42
1:AA:28:A:N3	1:AA:296:U:H4'	2.34	0.42
1:AA:44:A:O2'	1:AA:45:G:H5'	2.19	0.42
1:AA:493:A:H5'	1:AA:494:G:OP2	2.19	0.42
1:AA:537:G:H2'	1:AA:538:G:H8	1.85	0.42
1:AA:716:A:H2'	1:AA:717:U:C6	2.55	0.42
1:AA:86:G:N2	1:AA:87:C:H42	2.18	0.42
2:AC:25:THR:O	2:AC:28:PHE:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:117:VAL:HG12	3:AD:130:ASN:C	2.40	0.42
3:AD:162:GLU:C	3:AD:164:ARG:H	2.23	0.42
3:AD:90:LEU:CD2	3:AD:90:LEU:H	2.32	0.42
4:AE:19:ARG:O	4:AE:20:VAL:HB	2.19	0.42
1:AA:1114:C:H1'	21:AN:99:SER:O	2.20	0.42
20:AO:33:THR:HG23	20:AO:63:ARG:HH11	1.85	0.42
20:AO:36:ILE:HD11	20:AO:59:MET:HG3	2.02	0.42
13:AP:67:ILE:HD11	13:AP:71:VAL:CG2	2.45	0.42
14:AQ:47:ASP:C	14:AQ:49:ASN:H	2.23	0.42
14:AQ:59:GLU:O	14:AQ:75:VAL:HG22	2.20	0.42
16:AS:29:PRO:HD3	16:AS:47:THR:HB	2.02	0.42
17:AT:2:ASN:HD22	17:AT:3:ILE:HG13	1.84	0.42
22:BA:109:A:H2'	22:BA:110:C:H6	1.84	0.42
23:BB:1279:G:H2'	23:BB:1280:G:O4'	2.19	0.42
23:BB:1710:G:H4'	23:BB:2858:C:O2	2.18	0.42
23:BB:740:C:H5''	23:BB:1784:A:OP1	2.19	0.42
23:BB:1822:C:O2'	23:BB:1823:G:H5'	2.20	0.42
23:BB:2079:U:O2'	51:BZ:23:ASN:ND2	2.53	0.42
23:BB:2722:G:O2'	23:BB:2723:C:H5'	2.20	0.42
23:BB:532:A:C2'	23:BB:532:A:N3	2.80	0.42
23:BB:596:U:H2'	23:BB:597:G:C8	2.54	0.42
23:BB:686:U:H1'	36:B2:6:GLN:O	2.20	0.42
23:BB:724:U:H2'	23:BB:725:G:O4'	2.19	0.42
26:BD:73:VAL:O	26:BD:74:GLU:HB2	2.20	0.42
29:BE:15:SER:C	29:BE:17:THR:H	2.23	0.42
29:BE:145:ASP:CA	29:BE:166:LYS:HB3	2.50	0.42
48:BG:100:ASN:OD1	48:BG:101:VAL:HG13	2.19	0.42
40:BH:5:LEU:C	40:BH:7:ASP:H	2.22	0.42
41:BJ:133:ALA:C	41:BJ:135:GLN:H	2.23	0.42
41:BJ:3:THR:HB	41:BJ:44:TYR:HE1	1.83	0.42
34:B3:12:ARG:HG2	37:BL:62:PRO:O	2.20	0.42
38:BM:69:PRO:HG2	38:BM:70:ASP:H	1.85	0.42
28:BP:114:ASN:HD22	28:BP:114:ASN:HA	1.60	0.42
44:BQ:63:ARG:NH2	44:BQ:96:ASP:HA	2.34	0.42
49:BR:34:GLU:HB3	49:BR:58:VAL:CG2	2.50	0.42
45:BS:20:VAL:C	45:BS:22:ASP:N	2.72	0.42
50:BT:10:VAL:O	50:BT:12:ARG:N	2.51	0.42
35:BV:42:LEU:CD1	35:BV:47:VAL:HG21	2.49	0.42
52:BW:28:GLU:O	52:BW:30:VAL:N	2.52	0.42
52:BW:35:ILE:O	52:BW:36:ILE:C	2.58	0.42
1:CA:1019:A:H2'	1:CA:1020:G:C8	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1073:U:O2'	1:CA:1074:G:H5'	2.19	0.42
1:CA:977:A:N1	1:CA:1223:C:H2'	2.33	0.42
1:CA:1252:A:H2'	1:CA:1253:G:O4'	2.20	0.42
1:CA:1351:U:O2'	1:CA:1352:C:H5'	2.20	0.42
1:CA:456:A:H2'	1:CA:457:G:O4'	2.19	0.42
1:CA:833:G:O2'	1:CA:834:U:H5'	2.20	0.42
1:CA:921:U:H2'	1:CA:922:G:O4'	2.20	0.42
18:CB:15:PHE:CG	18:CB:16:GLY:N	2.87	0.42
18:CB:184:ALA:O	18:CB:199:ILE:HG13	2.19	0.42
2:CC:120:THR:HG21	2:CC:186:SER:HB2	2.02	0.42
4:CE:32:PHE:O	4:CE:51:LYS:HA	2.19	0.42
6:CG:147:ASN:C	6:CG:149:ALA:H	2.22	0.42
8:CI:16:ALA:HA	8:CI:65:THR:O	2.19	0.42
1:CA:779:C:H4'	10:CK:123:PRO:HA	2.01	0.42
21:CN:26:LEU:HD23	21:CN:31:SER:HB2	2.02	0.42
13:CP:24:SER:C	13:CP:25:ARG:HD3	2.40	0.42
13:CP:71:VAL:HA	13:CP:74:LEU:HG	2.01	0.42
16:CS:39:ILE:HG13	16:CS:39:ILE:H	1.36	0.42
33:D1:7:LYS:HZ3	34:D3:34:LYS:HG2	1.84	0.42
22:DA:28:C:H2'	22:DA:29:A:C1'	2.49	0.42
22:DA:52:A:H2'	22:DA:53:A:O4'	2.20	0.42
23:DB:1439:A:N7	23:DB:1440:U:C6	2.87	0.42
23:DB:1635:A:H2'	23:DB:1636:U:O4'	2.19	0.42
23:DB:1824:G:H2'	23:DB:1825:U:H6	1.85	0.42
23:DB:182:A:H2'	23:DB:183:C:C6	2.54	0.42
23:DB:1841:U:H2'	23:DB:1842:G:C8	2.54	0.42
23:DB:275:C:N3	23:DB:276:U:H1'	2.35	0.42
23:DB:728:G:O2'	23:DB:730:A:H8	2.02	0.42
23:DB:729:G:O4'	25:DC:206:LYS:NZ	2.48	0.42
23:DB:740:C:H5''	23:DB:1784:A:OP1	2.20	0.42
23:DB:898:C:O3'	23:DB:899:A:C4'	2.68	0.42
23:DB:974:G:H2'	23:DB:974:G:N3	2.35	0.42
25:DC:174:ARG:HG3	25:DC:180:MET:SD	2.60	0.42
26:DD:168:GLU:HG3	26:DD:170:VAL:CG1	2.48	0.42
29:DE:101:TYR:O	29:DE:104:ALA:HB3	2.19	0.42
47:DF:102:LEU:HD12	47:DF:103:ILE:HG13	2.01	0.42
47:DF:37:MET:HB2	47:DF:151:LEU:CB	2.50	0.42
22:DA:42:C:C1'	47:DF:65:LEU:HB2	2.50	0.42
40:DH:121:VAL:O	40:DH:122:LEU:HB2	2.19	0.42
40:DH:29:PHE:O	40:DH:30:LEU:C	2.57	0.42
23:DB:1138:G:H21	41:DJ:108:MET:HE1	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DN:72:ASP:O	42:DN:76:VAL:HG13	2.19	0.42
28:DP:31:VAL:CG1	28:DP:38:ARG:HG2	2.49	0.42
49:DR:34:GLU:HB3	49:DR:58:VAL:CG2	2.50	0.42
35:DV:42:LEU:CD1	35:DV:47:VAL:HG21	2.50	0.42
35:DV:80:HIS:CD2	35:DV:81:PRO:HD2	2.54	0.42
35:DV:35:GLU:HB3	35:DV:93:ARG:CZ	2.49	0.42
39:DX:8:GLU:HB3	39:DX:12:GLU:HB2	2.02	0.42
1:AA:1082:A:O2'	1:AA:1083:U:H5'	2.20	0.42
1:AA:1225:A:H3'	1:AA:1226:C:C5	2.55	0.42
1:AA:1293:C:O2'	1:AA:1294:G:H5'	2.19	0.42
1:AA:360:G:O2'	1:AA:361:G:H5'	2.20	0.42
1:AA:425:G:H2'	1:AA:426:U:C6	2.55	0.42
1:AA:467:U:H2'	1:AA:467:U:O2	2.20	0.42
1:AA:678:U:H2'	1:AA:679:C:H6	1.81	0.42
1:AA:934:C:H5''	56:AA:1724:HOH:O	2.20	0.42
18:AB:144:GLU:HG3	18:AB:148:GLY:HA3	2.02	0.42
18:AB:67:LEU:HD13	18:AB:68:PHE:N	2.35	0.42
2:AC:156:LEU:HD12	2:AC:163:ARG:HG3	2.02	0.42
2:AC:183:TYR:HA	2:AC:199:VAL:O	2.20	0.42
5:AF:1:MET:HB3	5:AF:66:ALA:C	2.40	0.42
5:AF:77:THR:O	5:AF:81:ASN:HB2	2.20	0.42
8:AI:14:SER:HA	8:AI:69:GLY:N	2.34	0.42
11:AL:42:LYS:HD2	11:AL:43:LYS:CG	2.49	0.42
16:AS:28:LYS:HB2	16:AS:29:PRO:HD2	2.01	0.42
19:AU:16:ARG:NH1	19:AU:19:LYS:NZ	2.68	0.42
31:B0:33:SER:O	31:B0:35:GLU:HG2	2.19	0.42
33:B1:3:GLY:C	33:B1:5:ARG:N	2.73	0.42
33:B1:8:ILE:HG13	33:B1:51:ALA:HA	2.02	0.42
23:BB:1262:A:H2'	23:BB:1263:U:O4'	2.20	0.42
23:BB:1463:C:H2'	23:BB:1464:G:H8	1.84	0.42
23:BB:1559:U:H3'	23:BB:1560:G:C5'	2.49	0.42
23:BB:1567:G:H1'	23:BB:1568:G:C5	2.55	0.42
23:BB:1684:G:H2'	23:BB:1685:C:C6	2.55	0.42
23:BB:1692:U:H2'	23:BB:1694:C:C4	2.54	0.42
23:BB:1729:U:H5'	23:BB:1730:C:OP2	2.20	0.42
23:BB:1998:A:O2'	23:BB:1999:C:H5'	2.20	0.42
23:BB:2074:U:H2'	23:BB:2075:U:C6	2.55	0.42
23:BB:215:G:C4'	23:BB:216:A:H4'	2.47	0.42
23:BB:2353:G:H1'	52:BW:30:VAL:CG1	2.50	0.42
23:BB:2454:G:H2'	23:BB:2455:G:H5'	2.00	0.42
23:BB:2466:C:OP1	32:B4:4:ARG:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1953:A:H2	23:BB:2549:G:N3	2.18	0.42
23:BB:2845:U:O2'	23:BB:2846:G:H5'	2.19	0.42
23:BB:265:A:N6	23:BB:427:U:O2'	2.52	0.42
23:BB:691:C:O2'	23:BB:692:C:H5'	2.19	0.42
26:BD:90:PHE:O	26:BD:91:THR:C	2.57	0.42
26:BD:34:VAL:HG13	26:BD:94:GLN:H	1.78	0.42
47:BF:62:GLN:HE21	47:BF:90:LEU:HD13	1.84	0.42
48:BG:10:VAL:HG23	48:BG:49:LEU:N	2.35	0.42
40:BH:81:ALA:CB	40:BH:147:VAL:H	2.28	0.42
40:BH:92:GLY:O	40:BH:123:ARG:HG2	2.20	0.42
41:BJ:1:MET:C	41:BJ:2:LYS:HZ3	2.23	0.42
41:BJ:58:ASN:O	41:BJ:59:ALA:HB3	2.20	0.42
37:BL:138:ALA:O	37:BL:139:GLY:C	2.58	0.42
38:BM:17:ASN:HD21	38:BM:95:LEU:CG	2.33	0.42
42:BN:119:SER:O	42:BN:120:GLU:HB2	2.19	0.42
50:BT:79:ASP:N	50:BT:79:ASP:OD2	2.53	0.42
39:BX:25:GLN:HB2	39:BX:46:VAL:HG11	2.02	0.42
39:BX:51:ALA:O	39:BX:52:ARG:C	2.58	0.42
39:BX:8:GLU:HB3	39:BX:12:GLU:HB2	2.01	0.42
1:CA:1110:A:O2'	1:CA:1111:A:H5'	2.19	0.42
1:CA:1163:A:O2'	1:CA:1164:G:H5'	2.19	0.42
1:CA:1220:G:H4'	16:CS:33:TRP:HB3	2.02	0.42
1:CA:1266:G:H22	1:CA:1268:G:H3'	1.85	0.42
1:CA:1332:A:O2'	1:CA:1333:A:H5'	2.19	0.42
1:CA:1288:A:H2	1:CA:1371:G:N3	2.18	0.42
1:CA:1372:U:C4	1:CA:1373:G:C5	3.08	0.42
1:CA:182:A:HO2'	1:CA:183:C:H3'	1.83	0.42
1:CA:222:C:H2'	1:CA:223:A:H8	1.85	0.42
1:CA:612:C:H2'	1:CA:613:C:C6	2.55	0.42
1:CA:833:G:H2'	1:CA:834:U:C6	2.54	0.42
18:CB:124:THR:HG23	18:CB:124:THR:O	2.18	0.42
19:AU:9:GLU:CG	2:CC:108:PRO:HG2	2.50	0.42
2:CC:19:SER:HA	2:CC:55:VAL:HG13	2.01	0.42
3:CD:202:LEU:O	3:CD:202:LEU:HD12	2.19	0.42
5:CF:11:HIS:CG	5:CF:12:PRO:HD2	2.54	0.42
6:CG:78:ARG:HH11	6:CG:78:ARG:C	2.23	0.42
8:CI:59:LYS:HZ3	8:CI:60:LEU:HG	1.85	0.42
9:CJ:40:ILE:HD12	9:CJ:73:LEU:HB3	2.01	0.42
13:CP:44:SER:C	13:CP:46:LYS:N	2.73	0.42
15:CR:28:LEU:C	15:CR:30:ASN:N	2.72	0.42
16:CS:15:LEU:O	16:CS:19:GLU:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CS:35:ARG:HG3	16:CS:36:ARG:N	2.35	0.42
17:CT:7:LYS:O	17:CT:11:ILE:HG13	2.20	0.42
19:CU:40:PRO:C	19:CU:42:THR:N	2.72	0.42
36:D2:13:ASN:O	36:D2:15:SER:N	2.52	0.42
34:D3:61:LEU:N	34:D3:62:PRO:CD	2.82	0.42
22:DA:13:G:O2'	22:DA:14:U:H3'	2.20	0.42
22:DA:80:U:H2'	22:DA:81:G:C8	2.55	0.42
23:DB:1392:A:N6	23:DB:1393:A:N6	2.68	0.42
23:DB:1404:C:H2'	23:DB:1405:U:H6	1.85	0.42
23:DB:1423:G:H2'	23:DB:1424:G:H8	1.84	0.42
23:DB:1535:A:O2'	23:DB:1537:G:N7	2.46	0.42
23:DB:1561:C:H2'	23:DB:1562:U:H6	1.84	0.42
23:DB:1695:G:H1'	25:DC:7:PRO:O	2.20	0.42
23:DB:1900:A:N1	23:DB:1970:A:C6	2.88	0.42
23:DB:2240:U:O2'	23:DB:2241:A:H5'	2.19	0.42
23:DB:2688:G:H1'	23:DB:2721:A:H61	1.84	0.42
23:DB:321:U:OP2	29:DE:130:LYS:HD3	2.19	0.42
23:DB:383:C:N3	23:DB:391:A:N6	2.68	0.42
25:DC:68:ARG:NH2	25:DC:190:THR:HG22	2.31	0.42
25:DC:230:PRO:O	25:DC:241:LYS:HD3	2.20	0.42
47:DF:127:TYR:CE2	47:DF:176:PHE:HE2	2.37	0.42
48:DG:8:VAL:HB	48:DG:49:LEU:O	2.20	0.42
40:DH:3:VAL:O	40:DH:18:GLN:HA	2.20	0.42
24:DI:54:ILE:O	24:DI:54:ILE:HG23	2.19	0.42
27:DK:88:ASN:C	27:DK:88:ASN:HD22	2.21	0.42
37:DL:90:VAL:HB	37:DL:122:VAL:HA	2.02	0.42
37:DL:132:ARG:CA	37:DL:135:ILE:HG22	2.49	0.42
37:DL:79:LEU:HB2	37:DL:113:ALA:N	2.31	0.42
49:DR:75:VAL:HG12	49:DR:76:LYS:N	2.34	0.42
45:DS:14:ALA:C	45:DS:16:LYS:N	2.73	0.42
50:DT:69:ARG:HA	50:DT:69:ARG:NE	2.27	0.42
52:DW:25:PHE:C	52:DW:27:GLY:N	2.71	0.42
51:DZ:14:THR:HA	51:DZ:28:ARG:HB2	2.02	0.42
1:AA:1455:G:H2'	1:AA:1456:A:H8	1.84	0.42
1:AA:153:C:H2'	1:AA:154:U:H6	1.84	0.42
1:AA:411:A:C4	1:AA:413:G:H1'	2.54	0.42
1:AA:818:G:H3'	1:AA:819:A:H5''	2.00	0.42
3:AD:97:LEU:O	3:AD:101:VAL:HG23	2.20	0.42
3:AD:106:PHE:CD1	3:AD:144:ILE:HD11	2.55	0.42
1:AA:410:G:OP1	3:AD:25:ARG:HD2	2.20	0.42
4:AE:156:ARG:HB2	4:AE:157:GLY:H	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:52:ARG:HH12	6:AG:121:ASN:HD22	1.66	0.42
7:AH:46:GLU:OE2	7:AH:63:LYS:HG3	2.20	0.42
7:AH:65:PHE:CG	7:AH:66:GLN:N	2.87	0.42
8:AI:5:TYR:O	8:AI:20:ILE:N	2.52	0.42
9:AJ:12:ALA:CB	9:AJ:18:ILE:HB	2.49	0.42
9:AJ:88:MET:SD	9:AJ:88:MET:N	2.92	0.42
10:AK:17:ASP:HB3	10:AK:80:ASN:OD1	2.20	0.42
21:AN:27:LYS:C	21:AN:29:ILE:N	2.73	0.42
16:AS:36:ARG:HE	16:AS:36:ARG:H	1.67	0.42
16:AS:66:VAL:HG23	16:AS:67:GLY:N	2.25	0.42
17:AT:4:LYS:HE3	17:AT:6:ALA:HB2	2.02	0.42
22:BA:75:G:N1	22:BA:102:G:N2	2.68	0.42
22:BA:42:C:C1'	47:BF:65:LEU:HB2	2.49	0.42
22:BA:8:C:O2'	43:BO:40:ILE:HD13	2.20	0.42
23:BB:1330:C:H2'	23:BB:1331:G:C8	2.55	0.42
23:BB:1338:G:O6	50:BT:66:LYS:HE2	2.19	0.42
23:BB:1366:A:H2'	23:BB:1367:A:O4'	2.19	0.42
23:BB:1518:C:H2'	23:BB:1519:G:H8	1.85	0.42
23:BB:1549:A:H2'	23:BB:1550:C:H6	1.82	0.42
23:BB:1930:G:C2'	23:BB:1931:U:OP2	2.68	0.42
23:BB:2076:U:O4'	23:BB:2076:U:O2	2.38	0.42
23:BB:2100:G:N3	23:BB:2100:G:H2'	2.35	0.42
23:BB:2144:G:O2'	23:BB:2145:C:H4'	2.20	0.42
23:BB:2271:G:H2'	23:BB:2272:U:C5	2.55	0.42
23:BB:2322:A:H2'	23:BB:2323:G:O4'	2.20	0.42
23:BB:2636:C:H2'	23:BB:2637:U:C6	2.55	0.42
23:BB:2654:A:H61	23:BB:2665:A:H3'	1.84	0.42
23:BB:2679:A:OP1	26:BD:114:LYS:HE2	2.20	0.42
23:BB:318:C:O2'	23:BB:319:G:H5'	2.20	0.42
23:BB:402:A:H2'	23:BB:403:U:H5'	2.02	0.42
23:BB:423:A:H5''	23:BB:424:G:C5'	2.49	0.42
23:BB:681:G:H2'	23:BB:682:G:H8	1.84	0.42
23:BB:709:U:H2'	23:BB:710:U:C6	2.54	0.42
23:BB:741:U:H2'	23:BB:742:A:C8	2.55	0.42
23:BB:798:G:O2'	23:BB:799:G:H5'	2.19	0.42
25:BC:159:THR:N	25:BC:194:VAL:CG1	2.83	0.42
26:BD:59:ARG:HB3	26:BD:59:ARG:NH2	2.35	0.42
29:BE:134:LEU:O	29:BE:138:LEU:HG	2.19	0.42
47:BF:33:ILE:CG2	47:BF:34:THR:N	2.82	0.42
47:BF:1:ALA:CB	47:BF:4:HIS:HB3	2.47	0.42
48:BG:148:ARG:HG2	48:BG:163:TYR:CZ	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BI:37:PHE:HZ	24:BI:56:VAL:HG11	1.85	0.42
24:BI:38:CYS:O	24:BI:42:ASN:ND2	2.53	0.42
24:BI:63:ASP:C	24:BI:65:SER:N	2.73	0.42
37:BL:80:SER:N	37:BL:113:ALA:HB3	2.32	0.42
38:BM:100:LYS:HD2	38:BM:101:VAL:H	1.85	0.42
42:BN:11:ASN:HA	42:BN:11:ASN:HD22	1.52	0.42
43:BO:49:VAL:HG23	43:BO:78:VAL:HG12	2.01	0.42
43:BO:93:ASP:C	43:BO:95:SER:H	2.24	0.42
28:BP:94:ALA:O	28:BP:95:LYS:HD2	2.19	0.42
45:BS:4:ILE:CG2	45:BS:106:VAL:HG22	2.49	0.42
46:BU:81:ARG:CD	46:BU:96:LYS:HD2	2.50	0.42
1:CA:1040:U:O2'	1:CA:1041:G:H5'	2.19	0.42
1:CA:1058:G:H2'	1:CA:1059:C:C6	2.55	0.42
1:CA:1112:C:N3	2:CC:177:LEU:HG	2.33	0.42
1:CA:1455:G:H2'	1:CA:1456:A:H8	1.85	0.42
1:CA:317:U:H2'	1:CA:318:G:H8	1.85	0.42
1:CA:373:A:O2'	1:CA:481:G:N2	2.52	0.42
1:CA:588:G:H2'	1:CA:588:G:N3	2.34	0.42
1:CA:862:C:O2'	1:CA:863:U:H5'	2.20	0.42
18:CB:107:ARG:HG3	18:CB:108:GLN:N	2.35	0.42
18:CB:53:LEU:HB2	18:CB:212:TYR:OH	2.20	0.42
18:CB:53:LEU:HA	18:CB:56:LEU:CD1	2.50	0.42
2:CC:46:LEU:CD1	2:CC:75:VAL:HG13	2.50	0.42
3:CD:1:ALA:O	3:CD:2:ARG:HG2	2.19	0.42
4:CE:39:GLY:HA3	4:CE:116:VAL:O	2.20	0.42
6:CG:105:GLU:HB3	6:CG:136:LYS:HZ3	1.83	0.42
6:CG:90:VAL:HG21	6:CG:95:ARG:HG3	2.01	0.42
7:CH:86:LYS:HG2	7:CH:124:ILE:HD11	2.01	0.42
8:CI:113:LYS:HG2	8:CI:114:LYS:N	2.34	0.42
1:CA:1129:C:C5'	8:CI:17:ARG:HH12	2.32	0.42
10:CK:125:LYS:O	19:CU:33:ARG:CZ	2.67	0.42
21:CN:67:GLY:O	21:CN:68:ARG:C	2.58	0.42
22:DA:106:G:H2'	22:DA:107:G:C8	2.55	0.42
22:DA:87:U:C2'	22:DA:88:C:O5'	2.68	0.42
23:DB:1301:A:N3	23:DB:1301:A:H2'	2.35	0.42
23:DB:1316:U:H2'	23:DB:1317:G:C8	2.55	0.42
23:DB:1494:A:H2'	23:DB:1495:A:H8	1.85	0.42
23:DB:1747:U:H2'	23:DB:1748:C:C6	2.54	0.42
23:DB:195:A:H3'	23:DB:196:A:H4'	2.01	0.42
23:DB:251:A:H2'	23:DB:252:G:O4'	2.19	0.42
23:DB:2708:G:H1'	42:DN:71:ARG:NH2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2807:U:H5'	23:DB:2808:G:OP2	2.20	0.42
23:DB:2811:G:O2'	23:DB:2812:G:H5'	2.20	0.42
23:DB:318:C:O2'	23:DB:319:G:H5'	2.20	0.42
23:DB:354:A:H2'	23:DB:355:U:C5	2.54	0.42
23:DB:479:A:H1'	23:DB:480:A:H5''	2.02	0.42
23:DB:63:A:O5'	23:DB:63:A:H8	2.03	0.42
23:DB:816:C:O2'	23:DB:817:C:H5'	2.20	0.42
23:DB:848:C:H2'	23:DB:849:A:H8	1.78	0.42
25:DC:219:VAL:HG12	25:DC:224:MET:HE2	2.02	0.42
23:DB:1657:U:O2'	26:DD:138:LEU:HD12	2.20	0.42
29:DE:105:LEU:HD21	29:DE:177:PRO:CB	2.50	0.42
47:DF:148:VAL:O	47:DF:149:ARG:CB	2.67	0.42
47:DF:76:PHE:HD2	47:DF:78:ILE:CD1	2.33	0.42
24:DI:5:GLN:O	24:DI:6:ALA:CB	2.68	0.42
41:DJ:17:VAL:HG22	41:DJ:55:ILE:HD11	2.02	0.42
27:DK:99:ILE:HG12	27:DK:115:ILE:HG13	2.01	0.42
38:DM:57:VAL:O	38:DM:59:ARG:N	2.47	0.42
42:DN:31:HIS:O	42:DN:32:GLU:HB2	2.20	0.42
43:DO:93:ASP:C	43:DO:95:SER:H	2.24	0.42
46:DU:5:ARG:HG2	46:DU:93:ARG:NH2	2.35	0.42
35:DV:14:LYS:O	35:DV:18:ARG:HB2	2.19	0.42
52:DW:30:VAL:HA	52:DW:60:ALA:O	2.19	0.42
39:DX:31:GLN:O	39:DX:36:GLN:HB2	2.19	0.42
51:DZ:56:MET:C	51:DZ:58:VAL:N	2.72	0.42
1:AA:104:G:O2'	1:AA:105:G:H5'	2.20	0.41
1:AA:1299:A:H5''	1:AA:1300:G:OP1	2.20	0.41
1:AA:1356:G:H3'	56:AA:1981:HOH:O	2.20	0.41
1:AA:190:A:C4	1:AA:191:G:H1'	2.54	0.41
1:AA:456:A:H2'	1:AA:457:G:O4'	2.20	0.41
1:AA:202:G:HO2'	1:AA:468:A:H8	1.66	0.41
1:AA:525:C:H2'	1:AA:526:C:H6	1.85	0.41
1:AA:600:A:H2'	1:AA:601:G:C8	2.55	0.41
1:AA:612:C:H2'	1:AA:613:C:C6	2.55	0.41
1:AA:875:U:H1'	7:AH:15:ASN:OD1	2.20	0.41
1:AA:926:G:H3'	1:AA:1505:G:N2	2.34	0.41
1:AA:982:U:OP2	21:AN:60:ARG:NH1	2.53	0.41
18:AB:156:LEU:HD12	18:AB:156:LEU:H	1.85	0.41
18:AB:17:HIS:CG	18:AB:18:GLN:H	2.38	0.41
3:AD:1:ALA:O	3:AD:2:ARG:HG2	2.19	0.41
3:AD:33:ILE:HD12	3:AD:34:GLU:H	1.85	0.41
4:AE:80:LEU:HG	4:AE:122:VAL:HG11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:70:PRO:HG3	6:AG:102:TRP:HH2	1.82	0.41
7:AH:63:LYS:HG2	7:AH:70:VAL:CG2	2.49	0.41
8:AI:79:ARG:O	8:AI:83:THR:HG22	2.20	0.41
11:AL:80:LEU:HB3	11:AL:97:VAL:HG23	2.01	0.41
14:AQ:80:LYS:H	14:AQ:80:LYS:HZ2	1.67	0.41
22:BA:13:G:N2	22:BA:16:G:N3	2.68	0.41
23:BB:1228:G:H2'	23:BB:1229:C:C6	2.54	0.41
23:BB:1286:A:C6	23:BB:1329:U:C2	3.08	0.41
23:BB:1316:U:O2'	23:BB:1317:G:H5'	2.19	0.41
23:BB:1427:A:H4'	23:BB:1428:C:O4'	2.20	0.41
23:BB:1435:G:H2'	23:BB:1436:G:C8	2.55	0.41
23:BB:1439:A:N7	23:BB:1440:U:C6	2.88	0.41
23:BB:1464:G:H2'	23:BB:1465:G:C8	2.55	0.41
23:BB:1418:G:H1'	23:BB:1580:A:N6	2.35	0.41
23:BB:20:C:O2'	23:BB:21:A:H5'	2.19	0.41
23:BB:223:A:N1	23:BB:407:G:O2'	2.50	0.41
23:BB:2324:U:H5'	23:BB:2325:G:H5''	2.01	0.41
23:BB:2479:U:H2'	23:BB:2480:C:H5'	2.02	0.41
23:BB:2860:A:H2'	23:BB:2861:U:O4'	2.20	0.41
23:BB:443:A:C5	29:BE:40:ARG:HD3	2.54	0.41
23:BB:605:G:O2'	23:BB:606:U:H5'	2.20	0.41
23:BB:637:A:OP2	37:BL:128:THR:HG21	2.20	0.41
23:BB:817:C:O2'	23:BB:839:U:H5''	2.20	0.41
23:BB:829:A:H5'	23:BB:831:G:N7	2.35	0.41
25:BC:180:MET:HB2	25:BC:268:ARG:CB	2.50	0.41
25:BC:89:ASN:HD22	25:BC:89:ASN:HA	1.64	0.41
26:BD:168:GLU:O	26:BD:169:ARG:C	2.57	0.41
48:BG:7:PRO:O	48:BG:8:VAL:CB	2.68	0.41
40:BH:116:ARG:HH22	40:BH:133:GLN:CB	2.33	0.41
41:BJ:29:ALA:O	41:BJ:33:ALA:N	2.46	0.41
41:BJ:45:THR:N	41:BJ:46:PRO:CD	2.83	0.41
41:BJ:77:HIS:ND1	41:BJ:79:GLY:N	2.67	0.41
27:BK:119:ALA:O	27:BK:120:PRO:C	2.58	0.41
23:BB:826:U:O2'	37:BL:53:GLY:HA3	2.20	0.41
38:BM:35:ALA:O	38:BM:36:VAL:HB	2.19	0.41
38:BM:17:ASN:HD21	38:BM:95:LEU:HG	1.85	0.41
42:BN:36:THR:OG1	42:BN:40:LYS:HD2	2.19	0.41
28:BP:99:LEU:HA	28:BP:102:ARG:HG3	2.02	0.41
27:BK:108:ARG:NH1	28:BP:34:GLY:HA2	2.33	0.41
28:BP:50:ARG:HB3	28:BP:57:ALA:O	2.20	0.41
28:BP:61:ARG:HD3	28:BP:70:GLU:OE1	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:BR:49:ILE:HG13	49:BR:49:ILE:H	1.59	0.41
45:BS:8:ARG:HA	45:BS:102:HIS:ND1	2.35	0.41
46:BU:46:LYS:HG2	46:BU:47:PRO:HD2	2.02	0.41
1:CA:1103:C:H6	1:CA:1103:C:O5'	2.02	0.41
1:CA:1325:C:O2'	1:CA:1326:U:H5'	2.19	0.41
1:CA:230:G:O2'	1:CA:231:U:H5'	2.20	0.41
1:CA:705:G:H2'	1:CA:706:A:H5'	2.02	0.41
1:CA:729:A:H2	1:CA:764:C:O2	2.02	0.41
18:CB:205:ALA:O	18:CB:209:VAL:HG13	2.20	0.41
2:CC:39:ARG:HG3	2:CC:39:ARG:NH1	2.35	0.41
7:CH:65:PHE:CG	7:CH:66:GLN:N	2.87	0.41
8:CI:103:VAL:CG2	8:CI:104:THR:N	2.82	0.41
1:CA:1151:A:C4'	9:CJ:41:PRO:HB2	2.50	0.41
11:CL:19:ASN:HA	11:CL:19:ASN:HD22	1.61	0.41
12:CM:64:VAL:C	12:CM:65:GLU:HG2	2.39	0.41
14:CQ:15:LYS:O	14:CQ:16:MET:C	2.57	0.41
15:CR:41:SER:HB3	15:CR:51:GLN:HB2	2.02	0.41
17:CT:66:ILE:HG23	17:CT:70:LYS:HD3	2.02	0.41
33:D1:35:LEU:O	33:D1:36:LYS:HB3	2.21	0.41
33:D1:39:ASP:OD1	33:D1:42:VAL:HG23	2.19	0.41
23:DB:591:U:H1'	34:D3:1:PRO:H2	1.85	0.41
34:D3:21:PHE:CZ	34:D3:58:ILE:HG23	2.55	0.41
23:DB:1322:A:O2'	23:DB:1323:C:H5'	2.19	0.41
23:DB:1628:G:O2'	23:DB:1629:U:H5'	2.20	0.41
23:DB:1670:C:H2'	23:DB:1670:C:O2	2.20	0.41
23:DB:2311:A:H3'	23:DB:2312:U:C6	2.55	0.41
23:DB:2379:G:H2'	23:DB:2380:C:C6	2.54	0.41
23:DB:26:G:H1'	23:DB:515:A:N6	2.35	0.41
23:DB:2700:A:H2'	23:DB:2701:U:H6	1.85	0.41
23:DB:2821:A:H2'	23:DB:2822:G:C8	2.55	0.41
23:DB:2853:C:H2'	23:DB:2854:G:C8	2.55	0.41
23:DB:285:G:H3'	23:DB:286:U:H6	1.84	0.41
23:DB:664:G:O2'	23:DB:665:U:H5'	2.20	0.41
23:DB:986:C:O2'	23:DB:987:C:H5'	2.20	0.41
25:DC:89:ASN:HA	25:DC:89:ASN:HD22	1.64	0.41
26:DD:102:ALA:C	26:DD:104:VAL:H	2.23	0.41
26:DD:118:PHE:N	26:DD:118:PHE:CD2	2.87	0.41
26:DD:32:ASN:HD22	26:DD:50:VAL:HG21	1.85	0.41
26:DD:35:THR:O	26:DD:71:ALA:HB1	2.20	0.41
26:DD:2:ILE:HG23	26:DD:84:LEU:HB3	2.02	0.41
29:DE:118:LEU:HD21	29:DE:188:MET:HE3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:134:GLN:HE21	47:DF:134:GLN:HB3	1.51	0.41
47:DF:101:ARG:HH11	47:DF:138:PRO:HB2	1.85	0.41
47:DF:137:PHE:CB	47:DF:138:PRO:HD2	2.43	0.41
38:DM:69:PRO:HG2	38:DM:70:ASP:H	1.84	0.41
38:DM:97:GLN:HB2	38:DM:98:PRO:HD2	2.01	0.41
42:DN:21:PHE:HA	42:DN:24:MET:HB3	2.02	0.41
43:DO:7:ARG:N	43:DO:10:ARG:HH11	2.18	0.41
50:DT:55:VAL:HG13	50:DT:85:VAL:HG12	2.02	0.41
51:DZ:64:ILE:CD1	51:DZ:64:ILE:H	2.23	0.41
1:AA:1071:C:H2'	1:AA:1072:G:H8	1.85	0.41
1:AA:1175:G:H4'	9:CJ:31:ARG:O	2.20	0.41
1:AA:1430:A:H2'	1:AA:1431:A:O4'	2.19	0.41
1:AA:160:A:H2'	1:AA:161:A:C8	2.55	0.41
1:AA:177:G:O4'	1:AA:177:G:N3	2.53	0.41
1:AA:16:A:C2'	1:AA:17:U:H5'	2.49	0.41
1:AA:193:C:H2'	1:AA:194:C:C5	2.55	0.41
1:AA:181:A:N6	1:AA:195:A:OP2	2.53	0.41
1:AA:295:C:H2'	1:AA:296:U:H6	1.85	0.41
1:AA:440:C:O2'	1:AA:441:A:H5'	2.20	0.41
1:AA:610:U:O2	1:AA:610:U:O4'	2.37	0.41
1:AA:647:C:H2'	1:AA:648:A:C8	2.55	0.41
1:AA:725:G:H2'	1:AA:726:C:H6	1.85	0.41
1:AA:764:C:N4	1:AA:812:G:N1	2.68	0.41
18:AB:112:ARG:NE	18:AB:116:LEU:HD11	2.32	0.41
18:AB:205:ALA:C	18:AB:207:ARG:H	2.23	0.41
3:AD:138:PRO:C	3:AD:140:ASP:H	2.23	0.41
1:AA:437:U:H4'	3:AD:151:GLN:HE21	1.85	0.41
3:AD:160:LEU:C	3:AD:162:GLU:N	2.73	0.41
3:AD:157:ALA:O	3:AD:160:LEU:HD22	2.20	0.41
3:AD:72:ARG:HA	3:AD:203:TYR:HE1	1.84	0.41
4:AE:84:VAL:HG11	4:AE:146:MET:HB3	2.00	0.41
6:AG:149:ALA:HB1	10:AK:58:THR:CB	2.46	0.41
10:AK:67:GLU:HG3	10:AK:68:ARG:N	2.34	0.41
12:AM:18:LEU:HD22	12:AM:32:ILE:HG21	2.02	0.41
12:AM:58:GLU:O	12:AM:61:LYS:HG2	2.20	0.41
14:AQ:80:LYS:HD2	14:AQ:80:LYS:C	2.40	0.41
17:AT:38:ILE:CD1	17:AT:82:ILE:HA	2.49	0.41
33:B1:8:ILE:HB	33:B1:24:LYS:HB2	2.01	0.41
22:BA:28:C:H2'	22:BA:29:A:C1'	2.50	0.41
23:BB:1597:A:C5'	23:BB:1598:A:H5'	2.35	0.41
23:BB:1668:A:O2'	23:BB:1674:G:N7	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:228:C:H4'	23:BB:229:C:C5'	2.49	0.41
23:BB:2350:C:O2'	23:BB:2351:G:H5'	2.19	0.41
23:BB:2471:A:O2'	23:BB:2472:G:C5'	2.68	0.41
23:BB:747:U:C4	23:BB:2613:U:C4	3.08	0.41
23:BB:2752:C:H2'	23:BB:2753:A:C8	2.55	0.41
23:BB:27:G:O5'	23:BB:27:G:H8	2.04	0.41
23:BB:303:G:H2'	23:BB:304:U:H6	1.83	0.41
23:BB:407:G:O2'	23:BB:408:G:H5'	2.19	0.41
23:BB:679:C:H2'	23:BB:680:C:C6	2.56	0.41
23:BB:928:A:O2'	30:BY:37:ARG:HD3	2.20	0.41
23:BB:984:A:O2'	23:BB:985:C:H5'	2.21	0.41
25:BC:155:ARG:HB3	25:BC:155:ARG:HH11	1.85	0.41
25:BC:204:LEU:HD23	25:BC:210:ALA:H	1.85	0.41
25:BC:239:PHE:O	25:BC:241:LYS:N	2.50	0.41
25:BC:51:ARG:O	25:BC:52:HIS:C	2.58	0.41
26:BD:172:VAL:O	26:BD:173:GLN:HB2	2.20	0.41
23:BB:674:G:H4'	29:BE:60:TRP:CZ2	2.55	0.41
47:BF:134:GLN:H	47:BF:150:GLY:CA	2.32	0.41
47:BF:137:PHE:N	47:BF:137:PHE:HD2	2.17	0.41
47:BF:38:GLY:HA2	47:BF:85:GLY:HA3	2.02	0.41
48:BG:17:LYS:HE3	48:BG:18:ILE:N	2.34	0.41
40:BH:54:LEU:O	40:BH:58:LEU:N	2.50	0.41
37:BL:79:LEU:N	37:BL:113:ALA:CB	2.83	0.41
23:BB:2277:G:P	38:BM:86:LYS:HB2	2.61	0.41
42:BN:87:PHE:HZ	42:BN:115:LEU:HB3	1.86	0.41
42:BN:47:VAL:O	42:BN:51:LEU:HD22	2.20	0.41
43:BO:115:LEU:CA	43:BO:116:GLN:HE21	2.33	0.41
44:BQ:60:TRP:CZ2	44:BQ:93:ILE:HB	2.55	0.41
49:BR:69:GLY:O	49:BR:90:ARG:HG3	2.21	0.41
35:BV:51:GLN:HA	35:BV:56:PHE:CD2	2.55	0.41
23:BB:2353:G:H1'	52:BW:30:VAL:HG12	2.02	0.41
52:BW:17:ALA:O	52:BW:36:ILE:HA	2.21	0.41
1:CA:1179:A:H2'	1:CA:1180:A:C8	2.55	0.41
1:CA:1426:G:H2'	1:CA:1427:C:H6	1.84	0.41
1:CA:1449:C:H2'	1:CA:1450:U:O4'	2.20	0.41
1:CA:411:A:C4	1:CA:413:G:H1'	2.55	0.41
1:CA:472:U:N3	1:CA:473:U:C4	2.88	0.41
1:CA:520:A:C2	1:CA:536:C:H1'	2.55	0.41
1:CA:5:U:H1'	1:CA:6:G:N1	2.34	0.41
1:CA:764:C:H2'	1:CA:765:G:C5'	2.47	0.41
1:CA:938:A:H2'	1:CA:939:G:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:968:A:H4'	1:CA:969:A:OP2	2.20	0.41
18:CB:111:LYS:HD3	18:CB:115:ASP:OD2	2.20	0.41
18:CB:24:PRO:C	18:CB:26:MET:H	2.23	0.41
18:CB:48:MET:CG	18:CB:200:PRO:HD2	2.49	0.41
1:CA:1056:U:H4'	2:CC:155:ARG:HG2	2.02	0.41
4:CE:104:ILE:O	4:CE:104:ILE:HG23	2.20	0.41
9:CJ:48:ARG:HG2	9:CJ:66:GLU:HB2	2.02	0.41
10:CK:12:ARG:HG2	10:CK:76:TYR:CE1	2.56	0.41
11:CL:35:ARG:NH1	11:CL:35:ARG:HA	2.34	0.41
15:CR:19:GLU:HG3	15:CR:54:LEU:HD12	2.01	0.41
16:CS:4:LEU:O	16:CS:6:LYS:N	2.53	0.41
10:CK:115:ILE:HG21	19:CU:23:GLU:OE1	2.20	0.41
31:D0:32:THR:OG1	31:D0:50:GLY:HA2	2.20	0.41
23:DB:2021:C:P	31:D0:8:THR:HG21	2.60	0.41
23:DB:1030:C:O2'	23:DB:1031:G:H5'	2.20	0.41
23:DB:1059:G:H2'	23:DB:1060:U:C5	2.55	0.41
23:DB:1764:C:O2'	23:DB:1765:U:H5'	2.19	0.41
23:DB:786:C:H5''	23:DB:1780:A:N7	2.36	0.41
23:DB:2060:A:C2	23:DB:2502:G:C5	3.07	0.41
23:DB:2064:C:H1'	23:DB:2450:A:C5	2.55	0.41
23:DB:222:A:H61	23:DB:232:G:H1'	1.85	0.41
23:DB:2385:C:H2'	23:DB:2386:A:C8	2.54	0.41
23:DB:277:G:HO2'	23:DB:361:G:H1	1.66	0.41
23:DB:265:A:N6	23:DB:427:U:O2'	2.53	0.41
23:DB:66:C:O2'	23:DB:67:U:H5'	2.20	0.41
23:DB:730:A:H5'	56:DB:3609:HOH:O	2.21	0.41
23:DB:941:A:H2	23:DB:1189:A:C2	2.37	0.41
26:DD:55:LYS:HZ3	26:DD:56:LYS:HG2	1.85	0.41
26:DD:91:THR:O	26:DD:92:VAL:C	2.58	0.41
48:DG:153:PRO:HB3	48:DG:158:GLY:CA	2.46	0.41
48:DG:10:VAL:HG23	48:DG:49:LEU:N	2.35	0.41
40:DH:38:PRO:HB2	40:DH:40:THR:HG23	2.02	0.41
24:DI:12:VAL:HG13	24:DI:41:PHE:CE2	2.55	0.41
23:DB:1063:G:C4'	24:DI:135:MET:HG2	2.50	0.41
38:DM:19:GLY:C	38:DM:20:LEU:HD22	2.40	0.41
38:DM:66:ARG:HE	38:DM:101:VAL:HG21	1.85	0.41
42:DN:33:ILE:CA	42:DN:114:GLU:HB2	2.49	0.41
31:D0:41:HIS:HB2	42:DN:99:LYS:O	2.20	0.41
44:DQ:65:ASN:O	44:DQ:69:ARG:N	2.47	0.41
45:DS:16:LYS:O	45:DS:19:LEU:HB3	2.20	0.41
45:DS:29:VAL:HG11	45:DS:55:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DS:71:VAL:O	45:DS:71:VAL:HG22	2.18	0.41
45:DS:86:MET:HB3	45:DS:94:ASP:HB2	2.02	0.41
50:DT:31:VAL:HG12	50:DT:84:TYR:CD2	2.54	0.41
50:DT:64:LYS:N	50:DT:64:LYS:HE3	2.34	0.41
46:DU:12:VAL:HG21	46:DU:38:ILE:HG12	2.02	0.41
35:DV:24:ASN:HB3	35:DV:44:HIS:HB3	2.01	0.41
35:DV:60:VAL:O	35:DV:61:LEU:HD23	2.19	0.41
52:DW:16:GLU:O	52:DW:17:ALA:C	2.58	0.41
23:DB:2352:A:N1	52:DW:30:VAL:CG1	2.83	0.41
52:DW:35:ILE:O	52:DW:36:ILE:C	2.59	0.41
51:DZ:70:GLU:O	51:DZ:71:LEU:C	2.57	0.41
1:AA:1015:G:O2'	1:AA:1016:A:H5'	2.21	0.41
1:AA:1121:U:O2'	1:AA:1122:U:H5'	2.20	0.41
1:AA:134:G:H2'	1:AA:135:C:O4'	2.20	0.41
1:AA:14:U:N3	1:AA:17:U:OP2	2.54	0.41
1:AA:512:U:O2'	1:AA:513:C:H5'	2.20	0.41
1:AA:652:U:H1'	1:AA:653:U:C5	2.56	0.41
1:AA:81:A:H3'	1:AA:83:C:C5	2.56	0.41
1:AA:929:G:H2'	1:AA:930:C:O4'	2.21	0.41
18:AB:18:GLN:HB2	18:AB:188:THR:OG1	2.20	0.41
18:AB:20:ARG:HG3	18:AB:20:ARG:HH11	1.85	0.41
2:AC:19:SER:CB	2:AC:21:TRP:HE1	2.32	0.41
2:AC:57:GLU:OE2	2:AC:64:ARG:HG2	2.20	0.41
2:AC:95:GLY:C	2:AC:96:VAL:HG22	2.41	0.41
3:AD:94:GLU:HA	3:AD:99:ASN:HD22	1.85	0.41
5:AF:47:LEU:HD21	5:AF:57:ALA:HB3	2.02	0.41
6:AG:53:SER:C	6:AG:55:LYS:N	2.74	0.41
8:AI:113:LYS:O	8:AI:113:LYS:HD3	2.19	0.41
10:AK:63:GLN:HG3	10:AK:98:ALA:CB	2.50	0.41
11:AL:106:VAL:CG2	11:AL:116:TYR:HB3	2.49	0.41
1:AA:1368:A:H5''	21:AN:100:TRP:HH2	1.86	0.41
20:AO:85:LEU:CB	20:AO:87:LEU:HG	2.49	0.41
14:AQ:30:HIS:HB3	14:AQ:34:GLY:N	2.35	0.41
15:AR:28:LEU:C	15:AR:30:ASN:N	2.72	0.41
1:AA:1319:A:OP1	16:AS:9:PHE:HB3	2.21	0.41
31:B0:28:SER:HB3	31:B0:37:HIS:HE1	1.85	0.41
32:B4:31:PRO:O	32:B4:34:LYS:HB3	2.20	0.41
23:BB:1417:C:C4'	23:BB:1587:G:H21	2.33	0.41
23:BB:1712:U:H3'	23:BB:1713:A:H2'	2.02	0.41
23:BB:1915:U:H2'	23:BB:1916:A:O4'	2.20	0.41
23:BB:201:C:H1'	23:BB:250:G:O6	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2752:C:H2'	23:BB:2753:A:H8	1.85	0.41
23:BB:2806:C:H2'	23:BB:2807:U:O4'	2.19	0.41
23:BB:294:A:H2'	23:BB:295:G:O5'	2.20	0.41
23:BB:523:C:O2'	23:BB:524:G:H5'	2.20	0.41
23:BB:63:A:H8	23:BB:63:A:P	2.43	0.41
23:BB:83:A:H5''	46:BU:1:ALA:N	2.35	0.41
25:BC:231:HIS:O	25:BC:232:GLY:C	2.58	0.41
29:BE:137:LYS:O	29:BE:141:MET:HG3	2.20	0.41
47:BF:84:ILE:CG2	47:BF:85:GLY:H	2.23	0.41
48:BG:93:TYR:HD2	48:BG:93:TYR:N	2.18	0.41
41:BJ:26:GLY:N	41:BJ:28:LEU:HD23	2.34	0.41
41:BJ:56:VAL:HG12	41:BJ:57:LEU:N	2.31	0.41
27:BK:18:ARG:HD3	27:BK:45:GLU:HG3	2.02	0.41
37:BL:90:VAL:HB	37:BL:122:VAL:HA	2.02	0.41
38:BM:57:VAL:O	38:BM:59:ARG:N	2.46	0.41
43:BO:83:LEU:CD1	43:BO:114:GLY:HA3	2.46	0.41
49:BR:14:VAL:HG22	49:BR:15:SER:H	1.85	0.41
46:BU:12:VAL:HG21	46:BU:38:ILE:CG1	2.50	0.41
1:CA:102:G:O2'	1:CA:103:U:H5'	2.20	0.41
1:CA:1032:G:H2'	1:CA:1033:G:C1'	2.51	0.41
1:CA:1178:G:H8	1:CA:1178:G:O5'	2.01	0.41
1:CA:1222:G:H2'	1:CA:1223:C:O4'	2.20	0.41
1:CA:1257:A:H3'	1:CA:1258:G:H5'	2.01	0.41
1:CA:1414:U:H2'	1:CA:1415:G:H8	1.84	0.41
1:CA:251:G:H22	1:CA:272:C:H41	1.69	0.41
1:CA:265:G:H2'	1:CA:267:C:H5	1.85	0.41
1:CA:467:U:O2	1:CA:467:U:H2'	2.20	0.41
1:CA:525:C:H2'	1:CA:526:C:H6	1.85	0.41
1:CA:694:A:C2	1:CA:695:A:H1'	2.55	0.41
1:CA:814:A:N7	1:CA:816:A:C4	2.88	0.41
18:CB:166:ASP:OD2	18:CB:190:SER:HA	2.21	0.41
2:CC:139:ASN:O	2:CC:143:LEU:HB3	2.21	0.41
2:CC:42:LEU:O	2:CC:46:LEU:HD23	2.20	0.41
3:CD:54:LEU:HD11	3:CD:55:ARG:HH21	1.83	0.41
3:CD:98:ASP:OD2	3:CD:99:ASN:N	2.53	0.41
5:CF:4:TYR:HA	5:CF:90:MET:O	2.20	0.41
5:CF:70:VAL:CG2	5:CF:71:ILE:N	2.83	0.41
7:CH:9:MET:HE1	7:CH:35:ILE:HB	2.03	0.41
8:CI:27:ILE:HD12	8:CI:27:ILE:H	1.85	0.41
14:CQ:34:GLY:O	14:CQ:35:LYS:C	2.58	0.41
15:CR:35:SER:O	15:CR:70:THR:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:CU:13:VAL:O	19:CU:13:VAL:HG13	2.20	0.41
36:D2:1:MET:CG	36:D2:2:LYS:H	2.32	0.41
23:DB:2420:C:OP1	34:D3:33:THR:HG22	2.20	0.41
22:DA:30:C:O2	22:DA:30:C:H2'	2.20	0.41
23:DB:1151:A:H2'	23:DB:1152:C:O4'	2.19	0.41
23:DB:1262:A:H2'	23:DB:1263:U:O4'	2.20	0.41
23:DB:1452:G:O2'	23:DB:1453:A:OP1	2.38	0.41
23:DB:1559:U:H3'	23:DB:1560:G:C5'	2.49	0.41
23:DB:1711:A:H2'	23:DB:1712:U:O4'	2.19	0.41
23:DB:1722:A:N6	23:DB:1738:G:H1'	2.34	0.41
23:DB:1860:G:O2'	23:DB:1861:G:H5'	2.20	0.41
23:DB:1892:C:O2'	23:DB:1893:C:H5'	2.20	0.41
23:DB:2000:C:O2'	23:DB:2001:C:H5'	2.20	0.41
23:DB:2100:G:C6	23:DB:2190:G:C5	3.09	0.41
23:DB:228:C:H4'	23:DB:229:C:C5'	2.48	0.41
23:DB:2365:G:H4'	52:DW:59:PHE:CE1	2.55	0.41
23:DB:2659:G:H1'	23:DB:2662:A:N6	2.34	0.41
23:DB:2756:U:C4	23:DB:2759:G:O6	2.73	0.41
23:DB:285:G:H3'	23:DB:286:U:C6	2.56	0.41
23:DB:632:A:H2'	23:DB:633:A:C8	2.55	0.41
23:DB:644:A:O2'	23:DB:645:C:H2'	2.19	0.41
23:DB:945:A:H3'	23:DB:946:C:H5''	2.02	0.41
23:DB:968:C:O2'	23:DB:969:G:H5'	2.20	0.41
26:DD:182:ALA:O	26:DD:184:ARG:HG2	2.20	0.41
29:DE:198:GLU:H	29:DE:198:GLU:HG2	1.59	0.41
29:DE:173:THR:HA	29:DE:199:MET:HE1	2.01	0.41
29:DE:31:VAL:HG21	29:DE:104:ALA:CB	2.50	0.41
48:DG:136:ASP:OD2	48:DG:139:VAL:HG23	2.20	0.41
48:DG:1:SER:N	48:DG:4:ALA:HB3	2.36	0.41
37:DL:115:GLU:N	37:DL:115:GLU:OE1	2.53	0.41
37:DL:120:VAL:HG12	37:DL:121:THR:N	2.34	0.41
37:DL:138:ALA:O	37:DL:139:GLY:C	2.58	0.41
38:DM:74:THR:HG21	38:DM:86:LYS:HE3	2.02	0.41
42:DN:116:VAL:HG13	42:DN:117:ASP:N	2.35	0.41
44:DQ:104:ALA:C	44:DQ:106:THR:H	2.23	0.41
44:DQ:4:LYS:HE3	44:DQ:8:ILE:HD11	2.03	0.41
49:DR:58:VAL:C	49:DR:59:ILE:HG13	2.41	0.41
45:DS:73:LYS:O	45:DS:106:VAL:N	2.53	0.41
35:DV:44:HIS:O	35:DV:46:LYS:N	2.53	0.41
35:DV:29:ILE:HG13	35:DV:88:HIS:HE1	1.85	0.41
1:AA:1408:A:H61	1:AA:1492:A:N6	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:220:G:O2'	1:AA:221:C:H5'	2.20	0.41
1:AA:317:U:H2'	1:AA:318:G:C8	2.55	0.41
1:AA:629:A:H2'	1:AA:630:A:O4'	2.20	0.41
1:AA:776:G:N2	1:AA:803:G:N7	2.68	0.41
1:AA:911:U:O2'	1:AA:912:C:H5'	2.20	0.41
1:AA:922:G:C6	1:AA:923:A:C6	3.09	0.41
1:AA:95:C:H2'	1:AA:96:U:H6	1.86	0.41
18:AB:110:ILE:C	18:AB:112:ARG:N	2.73	0.41
2:AC:112:ALA:HB1	2:AC:199:VAL:HG23	2.02	0.41
3:AD:117:VAL:HA	3:AD:122:ILE:HG12	2.00	0.41
5:AF:71:ILE:O	5:AF:74:LEU:HB3	2.21	0.41
5:AF:98:GLU:HG2	5:AF:99:ALA:N	2.35	0.41
6:AG:12:LEU:HD22	6:AG:13:PRO:HD2	2.02	0.41
10:AK:113:THR:CG2	10:AK:114:PRO:HD2	2.51	0.41
10:AK:91:GLY:O	10:AK:93:GLU:N	2.53	0.41
1:AA:1228:C:P	12:AM:106:ARG:HH12	2.42	0.41
21:AN:20:PHE:O	21:AN:22:LYS:N	2.53	0.41
20:AO:57:LEU:C	20:AO:57:LEU:HD13	2.40	0.41
13:AP:44:SER:C	13:AP:46:LYS:N	2.71	0.41
31:B0:18:HIS:ND1	31:B0:18:HIS:N	2.67	0.41
33:B1:10:LEU:HD23	33:B1:35:LEU:HD21	2.02	0.41
36:B2:4:THR:O	36:B2:5:PHE:HB2	2.20	0.41
32:B4:17:VAL:HG12	32:B4:18:LYS:N	2.36	0.41
23:BB:128:C:C2	23:BB:129:C:C5	3.09	0.41
23:BB:1459:G:H8	23:BB:1461:C:H5	1.68	0.41
23:BB:1647:U:H3'	23:BB:1647:U:OP1	2.20	0.41
23:BB:2418:A:H2'	23:BB:2419:U:O4'	2.20	0.41
23:BB:2581:G:H4'	23:BB:2582:G:C8	2.55	0.41
23:BB:2732:G:H3'	23:BB:2733:A:H5'	2.01	0.41
23:BB:379:G:N1	23:BB:396:G:C6	2.88	0.41
23:BB:780:G:H21	23:BB:783:A:H62	1.69	0.41
25:BC:29:PHE:CE2	25:BC:31:PRO:HG2	2.56	0.41
29:BE:118:LEU:HA	29:BE:186:VAL:CG1	2.50	0.41
47:BF:161:SER:HB3	47:BF:164:GLU:HG3	2.03	0.41
48:BG:126:THR:HB	48:BG:129:GLU:HB2	2.02	0.41
48:BG:34:ARG:N	48:BG:34:ARG:HH11	2.17	0.41
40:BH:58:LEU:HD12	40:BH:61:VAL:CG1	2.51	0.41
41:BJ:31:GLU:O	41:BJ:34:ARG:HB3	2.21	0.41
23:BB:2708:G:H1'	42:BN:71:ARG:NH2	2.35	0.41
43:BO:14:ALA:O	43:BO:18:LEU:HD13	2.20	0.41
28:BP:83:ILE:HD13	28:BP:83:ILE:C	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:BT:30:ILE:HG22	50:BT:87:LEU:HD21	2.02	0.41
39:BX:6:LEU:C	39:BX:8:GLU:H	2.22	0.41
39:BX:7:ARG:O	39:BX:9:LYS:HD3	2.20	0.41
1:CA:1061:G:H2'	1:CA:1062:U:O4'	2.20	0.41
1:CA:1088:G:H2'	1:CA:1089:G:O4'	2.21	0.41
1:CA:1354:U:O2'	1:CA:1355:G:H5'	2.20	0.41
1:CA:1530:G:H2'	1:CA:1531:A:C8	2.55	0.41
1:CA:401:C:H2'	1:CA:402:G:H8	1.85	0.41
1:CA:407:U:H2'	1:CA:408:A:C8	2.55	0.41
1:CA:493:A:H5'	1:CA:494:G:OP2	2.19	0.41
1:CA:631:C:C3'	1:CA:632:U:H5'	2.50	0.41
18:CB:85:SER:HB2	18:CB:88:GLN:HE22	1.85	0.41
3:CD:157:ALA:C	3:CD:159:GLU:N	2.73	0.41
3:CD:162:GLU:C	3:CD:164:ARG:H	2.22	0.41
8:CI:51:LEU:C	8:CI:53:LEU:H	2.23	0.41
12:CM:24:VAL:HG23	12:CM:24:VAL:O	2.19	0.41
14:CQ:25:GLU:N	14:CQ:40:THR:HG23	2.35	0.41
16:CS:51:HIS:ND1	16:CS:52:ASN:N	2.68	0.41
33:D1:8:ILE:HG13	33:D1:51:ALA:HA	2.02	0.41
23:DB:1158:C:H4'	30:DY:30:ARG:HH21	1.84	0.41
23:DB:1366:A:H2'	23:DB:1367:A:O4'	2.20	0.41
23:DB:1474:U:H2'	23:DB:1475:G:H5'	2.01	0.41
23:DB:1552:A:C2'	23:DB:1553:A:H5'	2.50	0.41
23:DB:1822:C:O2'	23:DB:1823:G:H5'	2.21	0.41
23:DB:2109:U:H2'	23:DB:2110:G:O4'	2.21	0.41
23:DB:2206:C:O2'	23:DB:2207:C:H5'	2.21	0.41
23:DB:2437:G:H2'	23:DB:2438:U:C6	2.55	0.41
23:DB:2479:U:C2'	23:DB:2480:C:H5'	2.50	0.41
23:DB:2547:A:H4'	27:DK:29:HIS:NE2	2.34	0.41
23:DB:547:A:H5'	23:DB:548:G:C5	2.55	0.41
23:DB:707:G:H2'	23:DB:708:G:O4'	2.19	0.41
23:DB:775:G:C6	23:DB:794:A:C8	3.08	0.41
23:DB:822:G:H2'	23:DB:823:C:C6	2.56	0.41
23:DB:967:U:H2'	23:DB:968:C:H6	1.85	0.41
25:DC:151:GLY:O	25:DC:152:GLN:HG3	2.20	0.41
25:DC:32:LEU:HD22	25:DC:63:ILE:HG21	2.01	0.41
26:DD:102:ALA:H	26:DD:104:VAL:CG2	2.32	0.41
26:DD:107:VAL:HG12	26:DD:108:ASP:N	2.36	0.41
26:DD:178:VAL:HG23	26:DD:189:VAL:O	2.20	0.41
26:DD:191:GLY:O	26:DD:192:ALA:HB3	2.21	0.41
47:DF:92:GLY:O	47:DF:95:MET:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DH:5:LEU:N	40:DH:17:ASP:O	2.53	0.41
41:DJ:14:ASP:O	41:DJ:53:TYR:N	2.53	0.41
41:DJ:28:LEU:CG	41:DJ:29:ALA:N	2.83	0.41
41:DJ:36:LEU:HD21	41:DJ:122:LEU:HD12	2.02	0.41
37:DL:129:LYS:HG2	37:DL:129:LYS:H	1.57	0.41
49:DR:59:ILE:HA	49:DR:101:ILE:H	1.84	0.41
49:DR:45:GLU:O	49:DR:46:GLU:C	2.59	0.41
50:DT:62:VAL:C	50:DT:64:LYS:HZ2	2.23	0.41
46:DU:73:ASN:O	46:DU:74:ALA:HB3	2.20	0.41
46:DU:81:ARG:CD	46:DU:96:LYS:HD2	2.50	0.41
52:DW:46:ALA:HB3	52:DW:79:ILE:O	2.20	0.41
51:DZ:32:ASN:O	51:DZ:33:LEU:O	2.39	0.41
51:DZ:70:GLU:O	51:DZ:73:ALA:N	2.46	0.41
1:AA:1158:C:O2	1:AA:1158:C:H3'	2.21	0.41
1:AA:168:G:O2'	1:AA:169:C:H5'	2.20	0.41
1:AA:184:G:H4'	1:AA:224:U:O3'	2.21	0.41
1:AA:707:U:H2'	1:AA:708:C:C6	2.55	0.41
1:AA:82:G:C2	1:AA:88:U:H1'	2.55	0.41
1:AA:662:U:O2'	1:AA:836:G:H5''	2.20	0.41
18:AB:150:ILE:O	18:AB:153:MET:HB3	2.20	0.41
18:AB:15:PHE:CD1	18:AB:16:GLY:N	2.89	0.41
18:AB:221:ARG:HG3	18:AB:222:GLU:HG2	2.01	0.41
4:AE:104:ILE:O	4:AE:104:ILE:HG23	2.20	0.41
4:AE:54:GLU:HG2	4:AE:56:PRO:HD2	2.03	0.41
5:AF:11:HIS:CG	5:AF:12:PRO:HD2	2.56	0.41
6:AG:71:THR:N	6:AG:141:HIS:HE1	2.09	0.41
7:AH:118:ALA:HB3	7:AH:120:LEU:CD2	2.50	0.41
1:AA:1320:C:C2	16:AS:35:ARG:HD3	2.56	0.41
19:AU:9:GLU:HG2	2:CC:109:GLU:HG3	2.03	0.41
22:BA:111:U:H2'	22:BA:112:G:C8	2.55	0.41
22:BA:46:A:H2'	22:BA:47:C:O4'	2.21	0.41
23:BB:1060:U:C1'	23:BB:1062:G:H5'	2.50	0.41
23:BB:1092:C:OP1	23:BB:2475:C:H4'	2.21	0.41
23:BB:1210:G:H5''	23:BB:1211:C:H3'	2.02	0.41
23:BB:1214:A:H2'	23:BB:1215:G:O4'	2.20	0.41
23:BB:1327:A:N6	23:BB:1328:A:C2	2.89	0.41
23:BB:1794:A:O2'	23:BB:1795:C:H5'	2.21	0.41
23:BB:2011:U:H2'	23:BB:2012:G:O4'	2.20	0.41
23:BB:212:G:H2'	23:BB:213:A:C8	2.56	0.41
23:BB:2508:G:H2'	23:BB:2509:G:H8	1.84	0.41
23:BB:2869:G:H2'	23:BB:2870:C:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:418:C:H2'	23:BB:419:U:H6	1.85	0.41
23:BB:526:A:N6	23:BB:2626:C:C4'	2.83	0.41
23:BB:714:U:C2	23:BB:716:A:H3'	2.55	0.41
23:BB:877:A:H2'	23:BB:899:A:C2	2.55	0.41
25:BC:57:HIS:ND1	25:BC:58:LYS:N	2.69	0.41
25:BC:61:TYR:HA	25:BC:85:ASN:HD21	1.85	0.41
26:BD:114:LYS:HD2	26:BD:116:LYS:HE3	2.01	0.41
26:BD:150:GLN:O	26:BD:153:GLY:N	2.53	0.41
26:BD:24:VAL:HG23	26:BD:189:VAL:C	2.41	0.41
29:BE:5:LEU:HA	29:BE:5:LEU:HD23	1.84	0.41
47:BF:133:GLU:HA	47:BF:150:GLY:HA2	2.01	0.41
47:BF:1:ALA:H2	47:BF:2:LYS:CE	2.34	0.41
47:BF:28:PRO:HB2	47:BF:168:LEU:CD2	2.42	0.41
48:BG:126:THR:HB	48:BG:129:GLU:CG	2.50	0.41
37:BL:29:LYS:C	37:BL:31:GLY:H	2.23	0.41
23:BB:2393:U:H5''	37:BL:62:PRO:HG3	2.03	0.41
43:BO:52:SER:OG	43:BO:54:VAL:HG12	2.20	0.41
1:AA:1441:A:C2	28:BP:113:LEU:HD22	2.56	0.41
44:BQ:42:GLY:HA3	49:BR:75:VAL:CG2	2.48	0.41
44:BQ:43:GLN:NE2	49:BR:77:PHE:HB3	2.36	0.41
46:BU:82:VAL:O	46:BU:94:PHE:O	2.37	0.41
1:CA:1328:C:O2'	1:CA:1329:A:H5'	2.21	0.41
1:CA:1342:C:H4'	8:CI:125:GLN:HB2	2.03	0.41
1:CA:1451:U:H5''	1:CA:1452:C:C5	2.56	0.41
1:CA:1472:U:O2'	1:CA:1473:G:H5'	2.21	0.41
1:CA:333:U:O2'	1:CA:334:C:H5'	2.21	0.41
1:CA:346:G:C2'	1:CA:347:G:H5'	2.51	0.41
1:CA:934:C:H5''	56:CA:1894:HOH:O	2.21	0.41
1:CA:437:U:C5'	3:CD:151:GLN:HE21	2.33	0.41
3:CD:25:ARG:HD3	3:CD:26:ALA:HB2	2.02	0.41
3:CD:97:LEU:O	3:CD:101:VAL:HG23	2.21	0.41
5:CF:38:ARG:O	5:CF:62:MET:O	2.39	0.41
6:CG:105:GLU:HB3	6:CG:136:LYS:HZ1	1.85	0.41
7:CH:16:GLY:C	7:CH:18:ALA:N	2.74	0.41
8:CI:51:LEU:C	8:CI:53:LEU:N	2.74	0.41
8:CI:29:ILE:CG2	8:CI:66:VAL:HG12	2.51	0.41
12:CM:64:VAL:HB	12:CM:65:GLU:OE2	2.21	0.41
20:CO:39:LEU:HD22	20:CO:59:MET:HE2	2.02	0.41
1:CA:719:C:H2'	15:CR:38:ILE:HD11	2.02	0.41
19:CU:40:PRO:O	19:CU:44:ARG:HB2	2.20	0.41
22:DA:87:U:H2'	22:DA:88:C:H5''	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1042:G:H2'	23:DB:1043:C:H6	1.84	0.41
23:DB:1228:G:H2'	23:DB:1229:C:H6	1.84	0.41
23:DB:1228:G:O2'	23:DB:1229:C:H5'	2.21	0.41
23:DB:1316:U:H2'	23:DB:1317:G:O4'	2.19	0.41
23:DB:1463:C:H2'	23:DB:1464:G:C8	2.55	0.41
23:DB:1486:U:H2'	23:DB:1487:U:C6	2.55	0.41
23:DB:1729:U:C5	23:DB:1731:G:N2	2.89	0.41
23:DB:1908:C:H2'	23:DB:1909:C:H6	1.85	0.41
23:DB:2135:A:H61	23:DB:2156:G:C2'	2.33	0.41
23:DB:2370:G:H2'	23:DB:2371:G:O4'	2.19	0.41
23:DB:23:G:C6	23:DB:518:G:C6	3.09	0.41
23:DB:475:C:C5	23:DB:476:G:N7	2.88	0.41
23:DB:63:A:H8	23:DB:63:A:P	2.43	0.41
23:DB:743:A:H2'	23:DB:744:U:C6	2.55	0.41
23:DB:833:A:H1'	37:DL:52:GLY:H	1.85	0.41
25:DC:23:LEU:HD12	25:DC:23:LEU:HA	1.84	0.41
25:DC:93:VAL:HG12	25:DC:101:ARG:C	2.40	0.41
26:DD:55:LYS:HZ3	26:DD:56:LYS:N	2.18	0.41
26:DD:62:LYS:CB	26:DD:63:PRO:HD3	2.46	0.41
29:DE:131:THR:O	29:DE:135:ALA:N	2.54	0.41
29:DE:141:MET:O	29:DE:143:LEU:HG	2.20	0.41
29:DE:188:MET:HG2	29:DE:193:VAL:CG2	2.51	0.41
47:DF:86:CYS:O	47:DF:88:VAL:HG23	2.21	0.41
47:DF:9:ASP:OD2	47:DF:9:ASP:N	2.53	0.41
48:DG:100:ASN:OD1	48:DG:101:VAL:HG13	2.21	0.41
48:DG:151:ARG:HG2	48:DG:160:GLY:HA2	2.03	0.41
23:DB:2529:G:C5'	48:DG:174:LYS:HG3	2.50	0.41
41:DJ:31:GLU:O	41:DJ:34:ARG:HB3	2.20	0.41
27:DK:3:GLN:CG	27:DK:4:GLU:N	2.83	0.41
28:DP:29:VAL:O	28:DP:40:GLN:N	2.53	0.41
28:DP:36:LYS:HG2	28:DP:36:LYS:H	1.69	0.41
28:DP:91:VAL:HG23	28:DP:92:ARG:N	2.36	0.41
23:DB:2848:G:N7	28:DP:94:ALA:HB2	2.36	0.41
44:DQ:65:ASN:HA	44:DQ:75:TYR:HB2	2.02	0.41
44:DQ:63:ARG:NH1	44:DQ:96:ASP:HB2	2.14	0.41
1:AA:1014:A:H2'	1:AA:1015:G:O4'	2.20	0.41
1:AA:1077:G:N1	1:AA:1080:A:OP2	2.53	0.41
1:AA:1157:A:H4'	1:AA:1158:C:O5'	2.21	0.41
1:AA:1216:A:O2'	1:AA:1217:C:H5'	2.21	0.41
1:AA:1191:A:H61	54:AA:1661:SCM:H11	1.86	0.41
1:AA:173:U:H6	1:AA:198:G:HO2'	1.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:284:C:H2'	1:AA:285:C:C6	2.56	0.41
1:AA:782:A:H2'	1:AA:783:C:O4'	2.20	0.41
18:AB:13:VAL:HB	18:AB:207:ARG:NE	2.27	0.41
18:AB:93:HIS:HA	18:AB:94:ARG:HH21	1.85	0.41
18:AB:94:ARG:N	18:AB:94:ARG:HE	2.18	0.41
2:AC:63:ILE:HD12	2:AC:91:ALA:HA	2.01	0.41
4:AE:148:SER:O	4:AE:151:MET:HB2	2.20	0.41
4:AE:32:PHE:O	4:AE:51:LYS:HA	2.21	0.41
6:AG:139:ASP:HA	6:AG:142:ARG:NH1	2.35	0.41
6:AG:4:ARG:HG3	6:AG:5:VAL:N	2.36	0.41
9:AJ:51:VAL:HG23	21:AN:80:ARG:CD	2.47	0.41
9:AJ:66:GLU:HB3	21:AN:98:ALA:HB2	2.02	0.41
9:AJ:92:LEU:HB2	9:AJ:93:ALA:H	1.64	0.41
11:AL:42:LYS:HB3	11:AL:43:LYS:H	1.45	0.41
1:AA:1228:C:H5''	12:AM:106:ARG:HH22	1.86	0.41
12:AM:44:ILE:N	12:AM:44:ILE:HD12	2.35	0.41
15:AR:25:ILE:HD12	15:AR:26:ALA:N	2.35	0.41
1:AA:719:C:H2'	15:AR:38:ILE:CD1	2.51	0.41
33:B1:26:LYS:HD2	33:B1:30:PRO:HA	2.01	0.41
33:B1:32:LYS:O	33:B1:34:GLU:HG3	2.21	0.41
36:B2:12:ARG:HG2	36:B2:44:VAL:HG11	2.03	0.41
22:BA:94:A:H2'	22:BA:95:U:O4'	2.20	0.41
23:BB:1109:C:H3'	23:BB:1110:G:C8	2.56	0.41
23:BB:1146:C:H2'	23:BB:1147:A:C8	2.56	0.41
23:BB:1740:G:H2'	23:BB:1741:C:H6	1.86	0.41
23:BB:1925:C:H2'	23:BB:1926:U:O4'	2.21	0.41
23:BB:1958:C:O2'	23:BB:1959:G:H5'	2.20	0.41
23:BB:2723:C:H4'	42:BN:1:MET:SD	2.60	0.41
23:BB:388:G:N7	23:BB:390:U:H2'	2.36	0.41
23:BB:409:G:H2'	23:BB:410:G:H8	1.85	0.41
23:BB:53:A:H2'	23:BB:54:G:O4'	2.20	0.41
23:BB:621:A:H2'	23:BB:622:G:O4'	2.21	0.41
25:BC:130:PRO:HA	25:BC:187:CYS:O	2.20	0.41
25:BC:239:PHE:HB3	25:BC:240:GLY:H	1.73	0.41
40:BH:31:VAL:O	40:BH:32:PRO:C	2.58	0.41
40:BH:43:ASN:O	40:BH:47:PHE:HB2	2.21	0.41
24:BI:14:ALA:CB	24:BI:50:LYS:HA	2.50	0.41
41:BJ:4:PHE:HB3	41:BJ:44:TYR:CE1	2.54	0.41
38:BM:97:GLN:HB2	38:BM:98:PRO:HD2	2.03	0.41
42:BN:16:HIS:O	42:BN:19:ALA:N	2.53	0.41
42:BN:65:LEU:HD11	42:BN:69:ARG:NH2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BO:49:VAL:O	43:BO:50:ALA:HB2	2.21	0.41
28:BP:8:GLU:HA	28:BP:54:LEU:HD22	2.03	0.41
44:BQ:59:LEU:O	44:BQ:59:LEU:HD22	2.20	0.41
46:BU:64:ILE:CG1	46:BU:65:GLN:N	2.82	0.41
46:BU:73:ASN:O	46:BU:74:ALA:HB3	2.20	0.41
30:BY:52:PHE:CG	30:BY:53:MET:N	2.89	0.41
1:CA:1053:G:C3'	1:CA:1054:C:H5'	2.50	0.41
1:CA:1342:C:H1'	8:CI:125:GLN:CD	2.40	0.41
1:CA:1464:U:H2'	1:CA:1465:A:C8	2.55	0.41
1:CA:1505:G:H4'	1:CA:1506:U:H5''	2.02	0.41
1:CA:1509:C:H2'	1:CA:1510:C:C6	2.55	0.41
1:CA:814:A:H5'	1:CA:1511:G:H4'	2.03	0.41
1:CA:356:A:N3	1:CA:368:U:O2'	2.46	0.41
1:CA:511:C:HO2'	1:CA:512:U:H5''	1.84	0.41
1:CA:598:U:H2'	1:CA:599:C:H6	1.83	0.41
1:CA:633:G:H2'	1:CA:634:C:C6	2.54	0.41
1:CA:952:U:O2'	1:CA:953:G:H5'	2.20	0.41
18:CB:127:LYS:HD2	18:CB:127:LYS:C	2.41	0.41
18:CB:17:HIS:CG	18:CB:18:GLN:N	2.89	0.41
2:CC:122:GLN:O	2:CC:127:VAL:HG13	2.20	0.41
3:CD:53:GLN:HG2	3:CD:198:LEU:CD2	2.50	0.41
5:CF:61:LEU:HB3	5:CF:62:MET:H	1.71	0.41
6:CG:145:GLU:N	6:CG:148:LYS:HB2	2.35	0.41
6:CG:92:PRO:HG2	6:CG:93:VAL:H	1.85	0.41
7:CH:25:THR:HG22	7:CH:26:MET:H	1.86	0.41
11:CL:107:LYS:HD2	11:CL:107:LYS:H	1.85	0.41
21:CN:65:GLN:N	21:CN:65:GLN:HE21	2.14	0.41
20:CO:70:LEU:CG	20:CO:78:TYR:HB2	2.50	0.41
14:CQ:30:HIS:HB3	14:CQ:34:GLY:N	2.36	0.41
34:D3:54:LEU:HG	34:D3:58:ILE:CD1	2.46	0.41
23:DB:1060:U:C1'	23:DB:1062:G:H5'	2.50	0.41
23:DB:1076:C:H2'	23:DB:1077:A:C8	2.56	0.41
23:DB:1316:U:C2	23:DB:1317:G:C8	3.09	0.41
23:DB:135:U:O2'	23:DB:136:G:H5'	2.21	0.41
23:DB:1567:G:H2'	25:DC:82:TYR:HE1	1.85	0.41
23:DB:1351:C:H4'	23:DB:1572:A:O4'	2.21	0.41
23:DB:1710:G:H4'	23:DB:2858:C:O2	2.20	0.41
23:DB:1266:G:H22	23:DB:2012:G:H2'	1.85	0.41
23:DB:2710:C:H2'	23:DB:2711:A:C8	2.56	0.41
23:DB:2821:A:OP2	23:DB:2821:A:H3'	2.21	0.41
23:DB:2886:A:H2'	23:DB:2887:A:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:372:G:O2'	23:DB:373:U:P	2.79	0.41
23:DB:97:C:H2'	23:DB:98:G:O4'	2.20	0.41
25:DC:5:CYS:HB2	25:DC:15:VAL:O	2.21	0.41
25:DC:52:HIS:O	25:DC:53:ILE:HB	2.21	0.41
25:DC:69:ASN:O	25:DC:70:LYS:C	2.59	0.41
26:DD:69:ALA:CA	26:DD:73:VAL:HB	2.49	0.41
23:DB:38:A:H1'	29:DE:43:THR:HB	2.02	0.41
47:DF:137:PHE:HD2	47:DF:137:PHE:N	2.18	0.41
48:DG:10:VAL:HG12	48:DG:10:VAL:O	2.20	0.41
41:DJ:45:THR:H	41:DJ:46:PRO:CD	2.28	0.41
27:DK:116:ILE:C	27:DK:118:LEU:H	2.24	0.41
23:DB:1952:A:C2	27:DK:22:ILE:HD12	2.56	0.41
37:DL:124:GLY:H	37:DL:143:GLU:CG	2.29	0.41
29:DE:29:HIS:CE1	37:DL:8:PRO:HG3	2.56	0.41
42:DN:35:LYS:HG2	42:DN:112:TYR:CE1	2.55	0.41
28:DP:80:VAL:CG1	28:DP:81:ASP:N	2.83	0.41
45:DS:99:ARG:HG2	45:DS:99:ARG:H	1.65	0.41
39:DX:19:LEU:O	39:DX:24:GLU:HB2	2.20	0.41
1:AA:1473:G:O2'	1:AA:1474:U:H5'	2.21	0.41
1:AA:1492:A:N3	23:BB:1913:A:N1	2.69	0.41
1:AA:384:G:O2'	1:AA:385:C:H5'	2.21	0.41
1:AA:626:G:H2'	1:AA:627:G:C8	2.56	0.41
1:AA:684:U:H2'	1:AA:685:G:O4'	2.21	0.41
1:AA:81:A:N6	1:AA:86:G:C6	2.89	0.41
18:AB:212:TYR:O	18:AB:215:ALA:HB3	2.21	0.41
2:AC:2:GLN:HB2	2:AC:2:GLN:HE21	1.57	0.41
2:AC:1:GLY:C	2:AC:2:GLN:HG3	2.41	0.41
3:AD:107:GLY:C	3:AD:157:ALA:HB1	2.41	0.41
8:AI:80:HIS:O	8:AI:83:THR:HG22	2.20	0.41
9:AJ:32:THR:CG2	9:AJ:83:THR:HG23	2.51	0.41
13:AP:45:GLU:C	13:AP:47:GLU:H	2.23	0.41
17:AT:2:ASN:CG	17:AT:3:ILE:H	2.23	0.41
17:AT:53:MET:HA	17:AT:56:ILE:CG1	2.51	0.41
10:AK:92:ARG:HH22	19:AU:19:LYS:HG2	1.82	0.41
23:BB:1429:G:O2'	23:BB:1430:G:H5'	2.20	0.41
23:BB:1609:A:O2'	23:BB:1610:A:H5''	2.21	0.41
23:BB:1657:U:C2'	23:BB:1658:C:H5'	2.49	0.41
23:BB:1670:C:O2	23:BB:1670:C:H2'	2.19	0.41
23:BB:1771:C:N4	23:BB:1772:A:H62	2.18	0.41
23:BB:1853:A:N1	23:BB:2087:G:H1'	2.36	0.41
23:BB:2146:C:H4'	23:BB:2148:G:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2291:U:OP1	23:BB:2380:C:O2'	2.39	0.41
23:BB:1129:A:N6	23:BB:2491:U:H5''	2.35	0.41
23:BB:2660:A:H2'	23:BB:2661:G:C8	2.56	0.41
23:BB:26:G:H1'	23:BB:515:A:N6	2.35	0.41
23:BB:2700:A:H2'	23:BB:2701:U:C6	2.55	0.41
23:BB:454:A:H3'	23:BB:455:C:C5'	2.48	0.41
23:BB:632:A:H2'	23:BB:633:A:C8	2.56	0.41
23:BB:65:U:N3	23:BB:66:C:C5	2.89	0.41
23:BB:707:G:H2'	23:BB:708:G:O4'	2.20	0.41
23:BB:715:A:H2'	23:BB:716:A:C8	2.56	0.41
23:BB:857:G:H2'	23:BB:858:G:H5'	1.99	0.41
23:BB:926:G:H2'	23:BB:927:A:H8	1.86	0.41
25:BC:207:ALA:O	25:BC:208:GLY:C	2.59	0.41
23:BB:1566:A:H5'	25:BC:213:ARG:NH1	2.35	0.41
26:BD:91:THR:O	26:BD:92:VAL:C	2.59	0.41
29:BE:11:ALA:O	29:BE:12:LEU:HB2	2.19	0.41
48:BG:84:LYS:HG3	48:BG:131:VAL:CA	2.50	0.41
24:BI:35:MET:C	24:BI:35:MET:SD	2.99	0.41
24:BI:91:LYS:O	24:BI:94:LYS:HB2	2.21	0.41
23:BB:636:G:C6	37:BL:111:ILE:HD12	2.56	0.41
38:BM:41:LEU:C	38:BM:43:ALA:H	2.23	0.41
49:BR:14:VAL:CG2	49:BR:18:GLN:HG3	2.51	0.41
49:BR:20:VAL:HG12	49:BR:21:ARG:N	2.36	0.41
50:BT:54:GLU:CB	50:BT:88:LYS:HB2	2.49	0.41
46:BU:64:ILE:HD12	46:BU:65:GLN:H	1.85	0.41
35:BV:80:HIS:CD2	35:BV:81:PRO:HD2	2.55	0.41
23:BB:855:G:N3	52:BW:23:LYS:HE3	2.36	0.41
51:BZ:40:VAL:CG2	51:BZ:43:GLU:HB3	2.46	0.41
51:BZ:64:ILE:CD1	51:BZ:64:ILE:H	2.26	0.41
1:CA:174:A:C2'	1:CA:175:C:H5'	2.51	0.41
1:CA:233:C:O2'	1:CA:234:C:H5'	2.20	0.41
1:CA:40:C:H2'	1:CA:41:G:O4'	2.21	0.41
1:CA:454:G:H2'	1:CA:455:G:H8	1.85	0.41
1:CA:652:U:H1'	1:CA:653:U:C5	2.55	0.41
1:CA:778:G:O2'	1:CA:779:C:H5'	2.21	0.41
1:CA:792:A:N3	1:CA:794:A:C5	2.89	0.41
1:CA:975:A:O2'	1:CA:976:G:OP2	2.32	0.41
18:CB:162:VAL:O	18:CB:184:ALA:HA	2.21	0.41
2:CC:149:LYS:CG	2:CC:168:ARG:HB2	2.48	0.41
3:CD:154:VAL:O	3:CD:158:LEU:HG	2.21	0.41
4:CE:84:VAL:HG11	4:CE:146:MET:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CE:20:VAL:O	4:CE:31:SER:HB2	2.21	0.41
4:CE:56:PRO:O	4:CE:57:ALA:C	2.59	0.41
5:CF:79:ARG:HD3	5:CF:79:ARG:HA	1.89	0.41
6:CG:147:ASN:C	6:CG:149:ALA:N	2.73	0.41
8:CI:40:ARG:H	8:CI:44:ARG:HD3	1.85	0.41
10:CK:70:ALA:C	10:CK:72:ALA:N	2.74	0.41
11:CL:35:ARG:HH21	11:CL:36:VAL:CG2	2.29	0.41
12:CM:43:LYS:CD	12:CM:43:LYS:H	2.34	0.41
12:CM:1:ALA:H3	12:CM:8:ILE:HG22	1.86	0.41
16:CS:62:THR:HB	16:CS:65:MET:HB3	2.02	0.41
23:DB:2391:G:P	34:D3:34:LYS:HZ3	2.43	0.41
32:D4:7:VAL:HG23	32:D4:35:GLN:CD	2.41	0.41
32:D4:36:ARG:O	32:D4:37:GLN:O	2.39	0.41
23:DB:1029:A:H8	23:DB:1029:A:O5'	2.04	0.41
23:DB:121:G:H2'	23:DB:122:G:H8	1.86	0.41
23:DB:1249:U:O4'	44:DQ:3:VAL:HG21	2.21	0.41
23:DB:1326:U:H2'	23:DB:1327:A:O4'	2.20	0.41
23:DB:1439:A:C5	23:DB:1552:A:N6	2.88	0.41
23:DB:1468:U:H2'	23:DB:1522:A:N6	2.34	0.41
23:DB:1538:G:H8	23:DB:1538:G:OP2	2.04	0.41
23:DB:1572:A:O2'	23:DB:1573:G:H5'	2.20	0.41
23:DB:1669:A:N3	23:DB:1669:A:H2'	2.35	0.41
23:DB:1788:C:H2'	23:DB:1789:A:H8	1.86	0.41
23:DB:1829:A:H3'	23:DB:1830:C:H6	1.84	0.41
23:DB:1925:C:H2'	23:DB:1926:U:O4'	2.20	0.41
23:DB:1937:A:C8	23:DB:1939:U:H2'	2.55	0.41
23:DB:1989:G:H2'	23:DB:1990:C:O4'	2.21	0.41
23:DB:2179:C:C2'	23:DB:2179:C:O2	2.68	0.41
23:DB:2223:G:H2'	23:DB:2224:G:H5'	2.03	0.41
23:DB:2233:U:H2'	23:DB:2234:G:C8	2.55	0.41
23:DB:2458:G:H1'	23:DB:2460:U:O4	2.20	0.41
23:DB:250:G:C6	23:DB:251:A:C6	3.08	0.41
23:DB:2597:G:C6	23:DB:2598:A:N6	2.89	0.41
23:DB:2721:A:H2'	23:DB:2722:G:H8	1.86	0.41
23:DB:445:C:H2'	23:DB:446:G:C8	2.56	0.41
23:DB:485:C:H2'	23:DB:486:C:H6	1.86	0.41
23:DB:825:A:O2'	37:DL:54:GLN:NE2	2.51	0.41
23:DB:940:G:H2'	23:DB:941:A:O4'	2.21	0.41
26:DD:106:LYS:O	26:DD:107:VAL:CB	2.69	0.41
23:DB:2512:C:O2'	26:DD:159:LYS:HE2	2.21	0.41
26:DD:125:TRP:CH2	26:DD:160:LYS:HG2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:172:VAL:O	26:DD:173:GLN:HB2	2.21	0.41
26:DD:55:LYS:HZ3	26:DD:56:LYS:CG	2.33	0.41
26:DD:48:ILE:HG22	26:DD:82:PHE:O	2.20	0.41
29:DE:108:ILE:HD11	29:DE:181:ILE:HB	2.02	0.41
29:DE:31:VAL:HG21	29:DE:104:ALA:HB2	2.02	0.41
47:DF:172:PHE:C	47:DF:174:PHE:H	2.24	0.41
47:DF:33:ILE:HD12	47:DF:33:ILE:N	2.36	0.41
47:DF:76:PHE:HB2	47:DF:77:LYS:H	1.67	0.41
47:DF:38:GLY:HA2	47:DF:85:GLY:HA3	2.03	0.41
48:DG:140:ILE:HD12	48:DG:141:GLY:N	2.36	0.41
48:DG:88:LEU:HD12	48:DG:88:LEU:C	2.41	0.41
40:DH:21:VAL:HG21	40:DH:30:LEU:HD11	2.02	0.41
24:DI:53:PRO:CG	24:DI:77:VAL:HG11	2.50	0.41
24:DI:99:LYS:HB2	24:DI:140:GLU:OE1	2.20	0.41
41:DJ:58:ASN:O	41:DJ:59:ALA:HB3	2.21	0.41
41:DJ:77:HIS:HD2	41:DJ:84:ILE:HB	1.85	0.41
27:DK:119:ALA:O	27:DK:120:PRO:O	2.38	0.41
27:DK:18:ARG:HD3	27:DK:45:GLU:HG3	2.02	0.41
38:DM:73:ILE:HG21	38:DM:91:TYR:CE2	2.55	0.41
42:DN:114:GLU:HG2	42:DN:115:LEU:O	2.21	0.41
28:DP:45:VAL:N	28:DP:60:VAL:HG13	2.35	0.41
28:DP:56:SER:O	28:DP:74:GLN:HA	2.20	0.41
49:DR:26:ASP:O	49:DR:27:ILE:HG12	2.20	0.41
49:DR:31:GLU:H	49:DR:63:VAL:HG22	1.86	0.41
40:DH:27:ARG:HG3	51:DZ:60:ASP:OD1	2.20	0.41
1:AA:1009:U:H2'	1:AA:1010:U:C5	2.55	0.41
1:AA:1138:G:N3	1:AA:1138:G:H3'	2.36	0.41
1:AA:1228:C:H2'	1:AA:1229:A:C8	2.56	0.41
1:AA:1251:A:H2'	1:AA:1252:A:C8	2.56	0.41
1:AA:213:G:C8	1:AA:214:C:C5	3.09	0.41
1:AA:214:C:H2'	1:AA:215:C:C6	2.56	0.41
1:AA:389:A:H2'	1:AA:389:A:N3	2.36	0.41
1:AA:411:A:N6	1:AA:413:G:H21	2.18	0.41
1:AA:705:G:H2'	1:AA:706:A:H5'	2.02	0.41
1:AA:750:C:O2	20:AO:23:GLY:HA3	2.20	0.41
1:AA:796:C:H2'	1:AA:797:C:H6	1.86	0.41
1:AA:821:G:O2'	1:AA:822:U:H5'	2.20	0.41
1:AA:825:A:H2'	1:AA:826:C:H6	1.85	0.41
1:AA:949:A:O2'	1:AA:950:U:H5'	2.20	0.41
2:AC:10:ARG:NH1	2:AC:10:ARG:HG3	2.36	0.41
3:AD:25:ARG:HE	3:AD:25:ARG:HB2	1.71	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AH:14:ARG:CG	7:AH:15:ASN:N	2.84	0.41
8:AI:43:ALA:HA	8:AI:46:VAL:CG1	2.49	0.41
10:AK:35:ASP:OD1	10:AK:37:GLN:HB2	2.20	0.41
12:AM:13:HIS:HB2	12:AM:16:ILE:HB	2.01	0.41
12:AM:84:CYS:C	12:AM:88:LEU:HG	2.40	0.41
21:AN:60:ARG:CG	21:AN:62:ARG:HE	2.30	0.41
14:AQ:58:VAL:HG12	14:AQ:77:VAL:HG13	2.03	0.41
15:AR:22:TYR:C	15:AR:24:ASP:H	2.23	0.41
16:AS:50:VAL:O	16:AS:56:HIS:HA	2.20	0.41
16:AS:44:ILE:CA	16:AS:61:VAL:HB	2.48	0.41
17:AT:3:ILE:O	17:AT:4:LYS:HB2	2.20	0.41
31:B0:42:ILE:HG12	42:BN:99:LYS:O	2.21	0.41
34:B3:50:SER:C	34:B3:52:GLY:N	2.74	0.41
32:B4:7:VAL:HG23	32:B4:35:GLN:CD	2.41	0.41
23:BB:1147:A:O2'	23:BB:1148:U:H5'	2.20	0.41
23:BB:1635:A:H2'	23:BB:1636:U:H5'	2.02	0.41
23:BB:1692:U:O2'	23:BB:1693:U:H2'	2.21	0.41
23:BB:1711:A:H2'	23:BB:1712:U:O4'	2.21	0.41
23:BB:1848:A:H2'	23:BB:1849:G:C8	2.55	0.41
23:BB:2249:U:H1'	23:BB:2275:C:H41	1.85	0.41
23:BB:2363:G:O2'	23:BB:2364:C:H5'	2.20	0.41
23:BB:2370:G:H2'	23:BB:2371:G:O4'	2.21	0.41
23:BB:2387:U:H1'	52:BW:38:ARG:CZ	2.50	0.41
23:BB:2573:C:H3'	56:BB:3614:HOH:O	2.21	0.41
23:BB:2660:A:H2'	23:BB:2661:G:O4'	2.21	0.41
23:BB:271:G:H1'	23:BB:272:A:C8	2.56	0.41
23:BB:353:C:O2	23:BB:353:C:H2'	2.21	0.41
23:BB:455:C:C3'	23:BB:456:C:H5'	2.51	0.41
23:BB:705:A:O2'	25:BC:6:LYS:HB2	2.20	0.41
25:BC:219:VAL:HG12	25:BC:224:MET:HE2	2.03	0.41
26:BD:154:LYS:HD3	26:BD:154:LYS:N	2.35	0.41
26:BD:38:LYS:HG3	26:BD:47:ALA:HB3	2.02	0.41
26:BD:3:GLY:O	26:BD:4:LEU:HD13	2.20	0.41
29:BE:138:LEU:O	29:BE:142:ALA:N	2.53	0.41
47:BF:121:PHE:CA	47:BF:127:TYR:HA	2.49	0.41
47:BF:92:GLY:O	47:BF:95:MET:HB3	2.20	0.41
48:BG:26:LYS:HA	48:BG:32:LEU:CA	2.51	0.41
48:BG:93:TYR:O	48:BG:94:ARG:HG3	2.21	0.41
40:BH:3:VAL:O	40:BH:18:GLN:HA	2.20	0.41
41:BJ:34:ARG:HH11	41:BJ:39:LYS:HG2	1.84	0.41
37:BL:111:ILE:CG2	37:BL:112:LEU:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BN:33:ILE:CB	42:BN:114:GLU:HB2	2.51	0.41
28:BP:80:VAL:CG1	28:BP:81:ASP:N	2.84	0.41
44:BQ:91:ARG:NH1	49:BR:11:GLN:H	2.19	0.41
44:BQ:86:SER:HB3	49:BR:51:VAL:HA	2.02	0.41
45:BS:14:ALA:C	45:BS:16:LYS:N	2.73	0.41
45:BS:1:MET:SD	45:BS:62:ASP:HB3	2.61	0.41
50:BT:25:GLU:OE1	50:BT:30:ILE:HA	2.21	0.41
50:BT:40:LYS:CA	50:BT:43:ILE:HG22	2.51	0.41
35:BV:2:PHE:CE2	35:BV:55:GLU:HB3	2.56	0.41
1:CA:1000:A:O2'	1:CA:1001:C:H5'	2.19	0.41
1:CA:1004:A:H2'	1:CA:1005:A:C8	2.56	0.41
1:CA:1320:C:P	16:CS:69:LYS:HG3	2.60	0.41
1:CA:177:G:N3	1:CA:177:G:O4'	2.53	0.41
1:CA:239:U:H5''	1:CA:239:U:H6	1.85	0.41
1:CA:321:A:H2'	1:CA:322:C:H6	1.86	0.41
1:CA:341:C:O2'	1:CA:342:C:H5'	2.20	0.41
1:CA:810:C:H2'	1:CA:811:C:O4'	2.20	0.41
1:CA:838:G:H2'	1:CA:839:C:H6	1.82	0.41
1:CA:874:G:O2'	1:CA:875:U:H5'	2.21	0.41
1:CA:964:A:H5''	56:CA:1951:HOH:O	2.20	0.41
4:CE:93:VAL:HA	4:CE:126:ALA:CB	2.50	0.41
4:CE:148:SER:HB2	4:CE:150:GLU:OE2	2.21	0.41
4:CE:91:SER:OG	4:CE:135:VAL:HG22	2.19	0.41
6:CG:74:VAL:HA	6:CG:86:VAL:O	2.21	0.41
7:CH:118:ALA:HB3	7:CH:120:LEU:CD2	2.51	0.41
4:CE:158:LYS:HZ2	7:CH:63:LYS:HD3	1.86	0.41
7:CH:76:ARG:HA	7:CH:76:ARG:HD2	1.91	0.41
21:CN:42:ASN:C	21:CN:44:VAL:H	2.23	0.41
20:CO:45:GLU:HB2	20:CO:46:HIS:ND1	2.35	0.41
1:CA:625:U:H4'	13:CP:16:PHE:CZ	2.55	0.41
15:CR:40:PRO:HB2	15:CR:43:ILE:HG12	2.03	0.41
1:CA:332:G:OP2	17:CT:2:ASN:HB3	2.20	0.41
17:CT:38:ILE:HD13	17:CT:38:ILE:HA	1.96	0.41
17:CT:53:MET:HA	17:CT:56:ILE:HG13	2.02	0.41
17:CT:85:LEU:CD2	17:CT:86:ALA:H	2.31	0.41
17:CT:85:LEU:HG	17:CT:86:ALA:N	2.36	0.41
32:D4:17:VAL:HG12	32:D4:18:LYS:N	2.36	0.41
22:DA:113:C:H2'	22:DA:114:C:H6	1.84	0.41
23:DB:1486:U:O2'	23:DB:1487:U:H5'	2.20	0.41
23:DB:1537:G:H5''	23:DB:1537:G:N3	2.36	0.41
23:DB:1609:A:O2'	23:DB:1610:A:H5''	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1409:C:C1'	23:DB:1913:A:H62	2.31	0.41
23:DB:1936:A:N6	23:DB:1963:U:N3	2.66	0.41
23:DB:2201:G:H2'	23:DB:2202:U:H6	1.86	0.41
23:DB:962:G:H21	23:DB:2250:G:H22	1.68	0.41
23:DB:2758:A:H2'	23:DB:2759:G:O4'	2.21	0.41
23:DB:332:A:O2'	23:DB:334:C:OP2	2.39	0.41
23:DB:362:A:H5'	23:DB:363:G:OP2	2.21	0.41
23:DB:396:G:O2'	23:DB:397:U:H5'	2.20	0.41
23:DB:715:A:H2'	23:DB:716:A:O4'	2.20	0.41
23:DB:724:U:H2'	23:DB:725:G:O4'	2.21	0.41
23:DB:878:A:H1'	23:DB:899:A:H62	1.85	0.41
23:DB:927:A:H2'	23:DB:928:A:C8	2.56	0.41
25:DC:196:ASN:O	25:DC:197:ALA:HB3	2.21	0.41
25:DC:231:HIS:O	25:DC:232:GLY:C	2.59	0.41
25:DC:51:ARG:O	25:DC:52:HIS:C	2.58	0.41
26:DD:16:THR:N	26:DD:20:VAL:O	2.54	0.41
29:DE:31:VAL:O	29:DE:34:ALA:HB3	2.21	0.41
47:DF:8:LYS:HD3	47:DF:9:ASP:N	2.36	0.41
24:DI:129:GLU:O	24:DI:133:ARG:HG3	2.21	0.41
27:DK:29:HIS:O	27:DK:30:ARG:C	2.58	0.41
27:DK:64:ARG:HA	27:DK:79:PHE:CE1	2.55	0.41
37:DL:2:ARG:HB3	37:DL:2:ARG:CZ	2.51	0.41
37:DL:79:LEU:HA	37:DL:79:LEU:HD23	1.87	0.41
37:DL:81:ASP:O	37:DL:82:LEU:HB2	2.21	0.41
42:DN:43:GLU:O	42:DN:46:ARG:HG2	2.21	0.41
42:DN:97:ILE:HG22	42:DN:113:ILE:HD12	2.02	0.41
46:DU:7:ASP:HA	46:DU:24:VAL:O	2.21	0.41
35:DV:53:LYS:NZ	35:DV:53:LYS:HA	2.36	0.41
52:DW:19:ARG:CD	52:DW:36:ILE:HD11	2.50	0.41
1:AA:1089:G:C6	1:AA:1090:U:C2	3.09	0.41
1:AA:1148:U:H5'	8:AI:6:TYR:HH	1.86	0.41
1:AA:1451:U:H5''	1:AA:1452:C:C5	2.55	0.41
1:AA:1472:U:O2'	1:AA:1473:G:H5'	2.21	0.41
1:AA:1417:G:H22	1:AA:1482:G:H2'	1.85	0.41
1:AA:222:C:H2'	1:AA:223:A:C8	2.56	0.41
1:AA:259:G:O2'	1:AA:260:G:H5'	2.21	0.41
1:AA:246:A:C2	1:AA:282:A:C5	3.09	0.41
1:AA:373:A:O2'	1:AA:481:G:N2	2.54	0.41
1:AA:471:U:H2'	1:AA:472:U:O4'	2.21	0.41
1:AA:778:G:H21	10:AK:121:ARG:HD3	1.86	0.41
1:AA:893:C:H2'	1:AA:894:G:H8	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:163:ARG:HG3	2:AC:163:ARG:O	2.21	0.41
2:AC:46:LEU:N	2:AC:46:LEU:HD22	2.36	0.41
3:AD:144:ILE:HG22	3:AD:145:ARG:H	1.86	0.41
6:AG:71:THR:HA	6:AG:90:VAL:CG2	2.42	0.41
7:AH:68:LYS:HG3	7:AH:69:ALA:O	2.20	0.41
8:AI:118:ARG:O	8:AI:118:ARG:HG2	2.21	0.41
1:AA:1350:A:OP1	8:AI:122:ARG:HD2	2.21	0.41
12:AM:105:ALA:C	12:AM:107:THR:H	2.25	0.41
33:B1:24:LYS:HG3	33:B1:25:ASN:H	1.86	0.41
32:B4:8:LYS:HB3	32:B4:8:LYS:HE2	1.91	0.41
23:BB:1079:C:H5	23:BB:1088:A:OP1	2.04	0.41
23:BB:1210:G:H5'	23:BB:1211:C:O5'	2.21	0.41
23:BB:1316:U:C2	23:BB:1317:G:C8	3.08	0.41
23:BB:1404:C:H2'	23:BB:1405:U:H6	1.86	0.41
23:BB:1458:U:H5'	23:BB:1460:U:OP2	2.19	0.41
23:BB:1914:C:H2'	23:BB:1915:U:C6	2.56	0.41
23:BB:2388:A:H5'	23:BB:2389:G:OP2	2.21	0.41
23:BB:2597:G:C6	23:BB:2598:A:N6	2.88	0.41
23:BB:275:C:H2'	23:BB:276:U:C5'	2.47	0.41
23:BB:2776:A:H4'	23:BB:2777:G:O5'	2.21	0.41
23:BB:2:G:H2'	23:BB:3:U:C6	2.56	0.41
23:BB:66:C:O2'	23:BB:67:U:H5'	2.21	0.41
23:BB:714:U:H3	23:BB:716:A:H3'	1.85	0.41
25:BC:54:GLY:O	25:BC:214:GLY:HA2	2.20	0.41
26:BD:102:ALA:H	26:BD:104:VAL:CG2	2.34	0.41
26:BD:107:VAL:HG12	26:BD:108:ASP:N	2.35	0.41
26:BD:8:LYS:HG2	26:BD:197:THR:N	2.34	0.41
23:BB:443:A:C4	29:BE:40:ARG:HD3	2.56	0.41
47:BF:33:ILE:HD12	47:BF:33:ILE:N	2.36	0.41
24:BI:83:ALA:N	24:BI:100:ILE:HD11	2.35	0.41
27:BK:39:ILE:HA	27:BK:39:ILE:HD13	1.90	0.41
37:BL:80:SER:HB3	37:BL:115:GLU:CD	2.41	0.41
42:BN:43:GLU:O	42:BN:46:ARG:HG2	2.21	0.41
28:BP:4:ILE:C	28:BP:6:GLN:N	2.74	0.41
28:BP:74:GLN:O	28:BP:75:THR:C	2.60	0.41
44:BQ:63:ARG:HH12	44:BQ:96:ASP:CB	2.17	0.41
45:BS:24:ILE:CG1	45:BS:36:LEU:HD21	2.51	0.41
39:BX:17:GLU:O	39:BX:20:ASN:HB2	2.21	0.41
1:CA:1059:C:O2'	1:CA:1060:U:H5'	2.21	0.41
1:CA:1132:C:H2'	1:CA:1133:G:H8	1.84	0.41
1:CA:1340:A:C2	1:CA:1341:U:H1'	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1444:U:H2'	1:CA:1445:U:C6	2.55	0.41
1:CA:384:G:O2'	1:CA:385:C:H5'	2.20	0.41
1:CA:433:G:H2'	1:CA:434:U:O4'	2.21	0.41
1:CA:707:U:H2'	1:CA:708:C:C6	2.56	0.41
1:CA:737:C:H2'	1:CA:738:C:C6	2.55	0.41
1:CA:796:C:H2'	1:CA:797:C:H6	1.86	0.41
1:CA:818:G:C2'	1:CA:819:A:H5''	2.51	0.41
1:CA:825:A:H2'	1:CA:826:C:H6	1.86	0.41
1:CA:842:U:H3'	1:CA:842:U:OP1	2.21	0.41
18:CB:107:ARG:HA	18:CB:110:ILE:HG13	2.02	0.41
18:CB:48:MET:HG3	18:CB:200:PRO:HD2	2.01	0.41
8:CI:56:MET:HA	8:CI:59:LYS:HB2	2.01	0.41
9:CJ:81:GLU:H	9:CJ:82:LYS:NZ	2.18	0.41
10:CK:45:THR:O	10:CK:46:ALA:C	2.59	0.41
1:CA:1229:A:H62	12:CM:103:THR:HG22	1.85	0.41
12:CM:30:LYS:NZ	12:CM:40:GLU:HG3	2.35	0.41
20:CO:47:LYS:C	20:CO:49:ASP:N	2.74	0.41
16:CS:13:HIS:ND1	16:CS:13:HIS:N	2.69	0.41
16:CS:45:GLY:CA	16:CS:60:PHE:HB3	2.48	0.41
17:CT:20:ASN:HA	17:CT:23:ARG:HG2	2.02	0.41
17:CT:3:ILE:O	17:CT:4:LYS:HB2	2.21	0.41
22:DA:7:G:O2'	22:DA:8:C:H5'	2.20	0.41
23:DB:1244:A:H4'	29:DE:29:HIS:HE1	1.86	0.41
23:DB:1392:A:C6	23:DB:1393:A:N6	2.89	0.41
23:DB:1660:G:H2'	23:DB:1661:G:H8	1.86	0.41
23:DB:2144:G:O2'	23:DB:2146:C:C5'	2.69	0.41
23:DB:215:G:C4'	23:DB:216:A:H4'	2.48	0.41
23:DB:2291:U:OP1	23:DB:2380:C:O2'	2.38	0.41
23:DB:2694:G:H2'	23:DB:2695:U:H6	1.85	0.41
23:DB:272:A:H2'	23:DB:273:G:C8	2.56	0.41
23:DB:1051:G:H4'	23:DB:2752:C:O2'	2.21	0.41
23:DB:2789:C:H2'	23:DB:2790:U:H5'	2.03	0.41
23:DB:629:G:H2'	23:DB:630:G:C8	2.55	0.41
23:DB:870:U:C3'	23:DB:871:U:H5''	2.51	0.41
23:DB:936:A:H2'	23:DB:937:C:H6	1.86	0.41
25:DC:106:PRO:HA	25:DC:195:GLY:N	2.36	0.41
26:DD:33:ARG:HH12	26:DD:53:GLY:C	2.24	0.41
26:DD:73:VAL:O	26:DD:74:GLU:HB2	2.20	0.41
26:DD:73:VAL:HG22	26:DD:74:GLU:N	2.36	0.41
29:DE:199:MET:HB3	29:DE:200:LEU:HD22	2.03	0.41
29:DE:69:ARG:O	29:DE:70:SER:CB	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:103:ILE:CG2	47:DF:173:ASP:HB2	2.51	0.41
48:DG:69:ALA:O	48:DG:73:SER:N	2.52	0.41
40:DH:126:GLY:H	40:DH:146:VAL:HB	1.86	0.41
41:DJ:88:THR:O	41:DJ:91:GLU:HB2	2.20	0.41
37:DL:29:LYS:C	37:DL:31:GLY:N	2.74	0.41
42:DN:82:GLU:C	42:DN:84:GLY:N	2.74	0.41
42:DN:55:ALA:HB1	42:DN:84:GLY:HA2	2.01	0.41
43:DO:15:ARG:HH21	43:DO:95:SER:HA	1.86	0.41
49:DR:78:ARG:HH21	49:DR:78:ARG:HG3	1.86	0.41
45:DS:1:MET:SD	45:DS:62:ASP:HB3	2.60	0.41
39:DX:30:MET:N	39:DX:30:MET:SD	2.94	0.41
39:DX:51:ALA:O	39:DX:52:ARG:C	2.58	0.41
1:AA:1139:G:H4'	1:AA:1140:C:OP1	2.21	0.41
1:AA:1215:G:H3'	1:AA:1215:G:OP2	2.20	0.41
1:AA:1392:G:H2'	1:AA:1393:U:H6	1.85	0.41
1:AA:195:A:H1'	1:AA:222:C:HO2'	1.85	0.41
1:AA:265:G:H2'	1:AA:267:C:H5	1.86	0.41
1:AA:35:G:O2'	1:AA:36:C:H5'	2.21	0.41
1:AA:439:U:H2'	1:AA:440:C:H6	1.86	0.41
1:AA:461:A:H3'	1:AA:462:G:C4'	2.50	0.41
1:AA:729:A:H2	1:AA:764:C:O2	2.03	0.41
4:AE:45:VAL:CG1	4:AE:116:VAL:HG23	2.51	0.41
4:AE:155:LYS:O	4:AE:158:LYS:HE3	2.21	0.41
4:AE:56:PRO:O	4:AE:59:ILE:HG13	2.21	0.41
5:AF:70:VAL:CG2	5:AF:71:ILE:N	2.84	0.41
6:AG:35:LYS:O	6:AG:38:ALA:HB3	2.21	0.41
7:AH:35:ILE:O	7:AH:39:LEU:HG	2.21	0.41
8:AI:103:VAL:HG23	8:AI:104:THR:H	1.85	0.41
8:AI:56:MET:O	8:AI:58:GLU:N	2.43	0.41
11:AL:120:ARG:HG2	11:AL:120:ARG:HH11	1.86	0.41
12:AM:21:ILE:HG21	12:AM:24:VAL:HG13	2.03	0.41
21:AN:97:LYS:HD3	21:AN:97:LYS:HA	1.86	0.41
16:AS:14:LEU:HD22	16:AS:34:SER:OG	2.21	0.41
31:B0:41:HIS:O	31:B0:42:ILE:O	2.39	0.41
22:BA:80:U:H2'	22:BA:81:G:C8	2.55	0.41
23:BB:120:U:H4'	23:BB:121:G:H5'	2.02	0.41
23:BB:120:U:H5''	23:BB:122:G:OP2	2.20	0.41
23:BB:1700:A:H2'	23:BB:1701:A:H5'	2.01	0.41
23:BB:1936:A:N6	23:BB:1963:U:N3	2.66	0.41
23:BB:2503:A:O2'	23:BB:2505:G:OP2	2.36	0.41
23:BB:2861:U:O2'	23:BB:2862:G:H5'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:391:A:N3	23:BB:391:A:H2'	2.36	0.41
23:BB:416:U:H2'	23:BB:417:C:H6	1.84	0.41
23:BB:727:A:H2	25:BC:8:THR:HG21	1.86	0.41
23:BB:812:C:H2'	23:BB:813:U:C6	2.55	0.41
23:BB:870:U:C3'	23:BB:871:U:H5''	2.51	0.41
26:BD:122:VAL:HA	26:BD:127:PHE:H	1.84	0.41
29:BE:141:MET:O	29:BE:143:LEU:HG	2.21	0.41
29:BE:12:LEU:CD1	29:BE:14:VAL:HG13	2.51	0.41
29:BE:2:GLU:OE1	29:BE:13:THR:N	2.54	0.41
29:BE:32:VAL:O	29:BE:36:ALA:N	2.49	0.41
29:BE:31:VAL:O	29:BE:34:ALA:HB3	2.21	0.41
47:BF:129:MET:SD	47:BF:129:MET:N	2.94	0.41
23:BB:2305:U:O2	47:BF:150:GLY:HA3	2.20	0.41
22:BA:42:C:C6	47:BF:65:LEU:HD22	2.56	0.41
23:BB:2311:A:O2'	47:BF:84:ILE:HD13	2.20	0.41
40:BH:117:LEU:CD1	40:BH:130:VAL:HG22	2.50	0.41
40:BH:116:ARG:HB3	40:BH:131:SER:N	2.36	0.41
40:BH:14:SER:O	40:BH:15:LEU:C	2.59	0.41
40:BH:5:LEU:HD11	40:BH:19:VAL:HG21	2.02	0.41
40:BH:94:ILE:HA	40:BH:98:ASP:OD1	2.20	0.41
40:BH:99:ILE:HG21	40:BH:130:VAL:CG2	2.50	0.41
24:BI:73:PRO:HA	24:BI:74:PRO:HD3	1.99	0.41
41:BJ:15:TRP:CD2	41:BJ:53:TYR:HB3	2.56	0.41
27:BK:19:VAL:O	27:BK:41:ILE:HD11	2.21	0.41
27:BK:64:ARG:HA	27:BK:79:PHE:CE1	2.56	0.41
38:BM:19:GLY:CA	38:BM:97:GLN:HB2	2.48	0.41
42:BN:33:ILE:CA	42:BN:114:GLU:HB2	2.49	0.41
42:BN:37:THR:OG1	42:BN:40:LYS:HE2	2.20	0.41
28:BP:91:VAL:HG23	28:BP:92:ARG:N	2.35	0.41
44:BQ:26:ALA:C	44:BQ:28:SER:N	2.73	0.41
50:BT:32:LEU:H	50:BT:83:ALA:CB	2.34	0.41
39:BX:20:ASN:O	39:BX:25:GLN:N	2.54	0.41
39:BX:59:GLU:OE2	39:BX:59:GLU:N	2.54	0.41
40:BH:28:ASN:HD21	51:BZ:36:HIS:CE1	2.39	0.41
1:CA:1065:U:O4	1:CA:1190:G:O4'	2.38	0.41
1:CA:511:C:O2'	1:CA:512:U:H5''	2.21	0.41
1:CA:552:U:H5'	11:CL:82:ARG:NH1	2.29	0.41
1:CA:724:G:O2'	1:CA:725:G:H5'	2.21	0.41
18:CB:53:LEU:CD1	18:CB:216:VAL:HA	2.51	0.41
18:CB:95:TRP:CH2	18:CB:99:MET:HG3	2.56	0.41
3:CD:157:ALA:O	3:CD:160:LEU:HD22	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:10:LEU:HD22	3:CD:62:ARG:NH1	2.35	0.41
4:CE:113:VAL:O	4:CE:116:VAL:N	2.53	0.41
5:CF:10:VAL:HG12	5:CF:11:HIS:N	2.36	0.41
8:CI:21:LYS:O	8:CI:61:ASP:N	2.54	0.41
9:CJ:36:VAL:CG1	9:CJ:76:ILE:HA	2.39	0.41
10:CK:35:ASP:OD1	10:CK:37:GLN:HB2	2.20	0.41
10:CK:91:GLY:C	10:CK:93:GLU:N	2.73	0.41
12:CM:59:VAL:O	12:CM:59:VAL:HG22	2.21	0.41
12:CM:95:PRO:CG	12:CM:101:THR:HG22	2.51	0.41
21:CN:79:SER:O	21:CN:83:VAL:HG23	2.21	0.41
20:CO:33:THR:CG2	20:CO:87:LEU:HD11	2.51	0.41
20:CO:64:ARG:CZ	20:CO:88:ARG:HD2	2.51	0.41
20:CO:84:ARG:C	20:CO:85:LEU:HD12	2.41	0.41
14:CQ:3:LYS:HE2	14:CQ:3:LYS:N	2.36	0.41
15:CR:31:TYR:CE1	15:CR:54:LEU:HD11	2.55	0.41
15:CR:58:ILE:O	15:CR:61:ALA:HB3	2.21	0.41
19:CU:29:ALA:HA	19:CU:32:ARG:CD	2.51	0.41
22:DA:3:C:H2'	22:DA:4:C:C6	2.56	0.41
23:DB:1063:G:O2'	23:DB:1064:C:H5'	2.20	0.41
23:DB:1171:G:H2'	23:DB:1172:C:C6	2.56	0.41
23:DB:1317:G:H2'	23:DB:1318:U:O4'	2.21	0.41
23:DB:1439:A:C8	23:DB:1440:U:C6	3.09	0.41
23:DB:1418:G:H1'	23:DB:1580:A:N6	2.36	0.41
23:DB:1842:G:O2'	23:DB:1843:C:H5'	2.20	0.41
23:DB:2520:C:O2'	23:DB:2521:C:H5'	2.20	0.41
23:DB:2572:A:OP1	23:DB:2574:G:H4'	2.21	0.41
23:DB:2772:C:H2'	23:DB:2773:C:C6	2.56	0.41
23:DB:2812:G:H2'	23:DB:2813:A:C8	2.56	0.41
23:DB:296:U:O2'	23:DB:297:G:H5'	2.20	0.41
23:DB:419:U:H2'	23:DB:420:C:H6	1.84	0.41
23:DB:640:C:H2'	23:DB:641:U:C6	2.56	0.41
23:DB:826:U:H2'	23:DB:828:U:O4'	2.21	0.41
26:DD:117:GLY:HA2	26:DD:164:GLN:HE22	1.83	0.41
26:DD:154:LYS:HD3	26:DD:154:LYS:N	2.36	0.41
29:DE:11:ALA:O	29:DE:12:LEU:HB2	2.21	0.41
29:DE:17:THR:C	29:DE:19:PHE:N	2.74	0.41
23:DB:2312:U:C4'	47:DF:84:ILE:HG21	2.51	0.41
40:DH:4:ILE:HD13	40:DH:37:VAL:HG13	2.03	0.41
40:DH:41:LYS:HA	40:DH:44:ILE:HG12	2.02	0.41
41:DJ:26:GLY:C	41:DJ:28:LEU:N	2.74	0.41
41:DJ:26:GLY:N	41:DJ:28:LEU:HD23	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DK:67:LYS:HD3	27:DK:67:LYS:C	2.40	0.41
29:DE:29:HIS:CD2	37:DL:8:PRO:HA	2.55	0.41
42:DN:118:ARG:HG2	42:DN:118:ARG:O	2.21	0.41
42:DN:8:ARG:HG2	42:DN:10:LEU:HD22	2.02	0.41
43:DO:115:LEU:HD22	43:DO:115:LEU:N	2.36	0.41
28:DP:52:ARG:HD2	28:DP:52:ARG:HA	1.87	0.41
28:DP:83:ILE:C	28:DP:83:ILE:HD13	2.41	0.41
28:DP:94:ALA:O	28:DP:95:LYS:HD2	2.21	0.41
44:DQ:48:ASP:N	44:DQ:48:ASP:OD2	2.53	0.41
50:DT:59:ASN:O	50:DT:83:ALA:O	2.38	0.41
46:DU:5:ARG:HB2	46:DU:8:ASP:OD2	2.21	0.41
46:DU:64:ILE:CG1	46:DU:65:GLN:N	2.84	0.41
35:DV:38:LEU:CD1	35:DV:40:ILE:HG23	2.51	0.41
35:DV:83:LYS:HA	35:DV:84:PRO:HD3	1.91	0.41
52:DW:35:ILE:CG1	52:DW:35:ILE:O	2.69	0.41
51:DZ:5:CYS:HB3	51:DZ:10:LYS:H	1.85	0.41
1:AA:1256:A:N1	1:AA:1278:G:O2'	2.49	0.41
1:AA:971:G:N2	1:AA:1363:A:H8	2.19	0.41
1:AA:1367:C:O2'	1:AA:1368:A:H5'	2.20	0.41
1:AA:1375:A:O2'	1:AA:1376:U:H5'	2.21	0.41
1:AA:198:G:N7	1:AA:220:G:N2	2.68	0.41
1:AA:643:C:H2'	1:AA:644:U:C6	2.55	0.41
1:AA:708:C:O2'	1:AA:709:U:H5'	2.21	0.41
1:AA:730:G:O2'	1:AA:766:A:H5'	2.21	0.41
1:AA:796:C:H5'	10:AK:128:VAL:HG13	2.01	0.41
1:AA:814:A:H5'	1:AA:1511:G:H4'	2.02	0.41
1:AA:838:G:H2'	1:AA:839:C:H6	1.82	0.41
1:AA:968:A:H5'	1:AA:968:A:N3	2.35	0.41
1:AA:981:U:H4'	21:AN:60:ARG:HB3	2.03	0.41
1:AA:993:G:H2'	1:AA:995:C:H41	1.84	0.41
18:AB:204:ASP:O	18:AB:205:ALA:HB3	2.21	0.41
2:AC:179:ALA:HA	2:AC:205:GLU:CA	2.47	0.41
2:AC:52:SER:HB3	2:AC:114:LEU:CD2	2.49	0.41
2:AC:91:ALA:HB1	2:AC:96:VAL:O	2.20	0.41
2:AC:9:ILE:HG23	2:AC:10:ARG:N	2.36	0.41
4:AE:113:VAL:O	4:AE:116:VAL:N	2.54	0.41
4:AE:82:HIS:CE1	4:AE:146:MET:HA	2.55	0.41
6:AG:129:ASN:CA	6:AG:134:VAL:HG21	2.51	0.41
6:AG:16:LYS:HD3	6:AG:17:PHE:CE1	2.56	0.41
7:AH:43:GLY:HA2	7:AH:63:LYS:HZ1	1.86	0.41
9:AJ:37:ARG:HA	9:AJ:37:ARG:HE	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:59:GLY:O	11:AL:60:PHE:HD1	2.04	0.41
21:AN:60:ARG:NE	21:AN:62:ARG:HG2	2.36	0.41
21:AN:60:ARG:O	21:AN:61:ASN:HB2	2.20	0.41
20:AO:28:GLN:O	20:AO:32:LEU:HD23	2.21	0.41
20:AO:36:ILE:HD12	20:AO:60:VAL:HG22	2.02	0.41
14:AQ:43:LEU:N	14:AQ:43:LEU:HD12	2.35	0.41
17:AT:20:ASN:HA	17:AT:23:ARG:HG2	2.03	0.41
17:AT:59:ARG:HB2	17:AT:59:ARG:HH11	1.86	0.41
34:B3:30:HIS:CD2	34:B3:31:ILE:N	2.89	0.41
22:BA:13:G:O2'	22:BA:14:U:H3'	2.21	0.41
22:BA:39:A:O2'	22:BA:40:U:H5'	2.20	0.41
23:BB:1197:G:H2'	23:BB:1198:U:H6	1.85	0.41
23:BB:1526:C:H2'	23:BB:1527:G:C8	2.56	0.41
23:BB:1745:A:H2'	23:BB:1746:A:H8	1.85	0.41
23:BB:1841:U:C2	23:BB:1842:G:C8	3.09	0.41
23:BB:1936:A:H2	23:BB:1943:U:O4	2.04	0.41
23:BB:1956:U:O2	23:BB:1985:C:H4'	2.21	0.41
23:BB:2210:U:C4	23:BB:2212:A:N7	2.89	0.41
23:BB:2306:C:H6	23:BB:2306:C:O5'	2.04	0.41
23:BB:2428:G:H5''	23:BB:2429:G:H5'	2.03	0.41
23:BB:2804:U:H2'	23:BB:2805:C:H6	1.86	0.41
23:BB:422:A:H2'	23:BB:422:A:N3	2.36	0.41
23:BB:734:A:C2	23:BB:735:A:H1'	2.55	0.41
23:BB:822:G:H2'	23:BB:823:C:C6	2.55	0.41
25:BC:93:VAL:HG12	25:BC:101:ARG:C	2.41	0.41
26:BD:182:ALA:O	26:BD:184:ARG:HG2	2.19	0.41
29:BE:152:GLU:O	29:BE:153:LEU:HB3	2.21	0.41
47:BF:28:PRO:HA	47:BF:158:THR:OG1	2.20	0.41
47:BF:47:LYS:HA	47:BF:50:ASP:OD1	2.21	0.41
40:BH:133:GLN:HG3	40:BH:139:PHE:HD2	1.85	0.41
41:BJ:44:TYR:C	41:BJ:44:TYR:CD2	2.93	0.41
23:BB:1652:A:N6	42:BN:11:ASN:HD21	2.05	0.41
43:BO:30:ARG:HG2	43:BO:31:THR:H	1.85	0.41
43:BO:7:ARG:HA	43:BO:10:ARG:CZ	2.51	0.41
28:BP:6:GLN:HA	28:BP:9:GLN:CD	2.41	0.41
49:BR:60:LYS:O	49:BR:98:ILE:HA	2.20	0.41
45:BS:57:ASN:ND2	45:BS:57:ASN:N	2.69	0.41
50:BT:22:THR:HA	50:BT:25:GLU:HB3	2.03	0.41
35:BV:53:LYS:HE3	35:BV:55:GLU:OE1	2.20	0.41
23:BB:2230:G:H4'	51:BZ:31:PRO:O	2.20	0.41
51:BZ:5:CYS:C	51:BZ:7:VAL:H	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1321:U:H4'	12:CM:85:TYR:CE1	2.55	0.41
1:CA:1459:G:H2'	1:CA:1460:C:H6	1.85	0.41
1:CA:181:A:N6	1:CA:195:A:OP2	2.53	0.41
1:CA:202:G:HO2'	1:CA:468:A:H8	1.65	0.41
1:CA:20:U:O2'	1:CA:21:G:H5'	2.21	0.41
1:CA:250:A:H1'	1:CA:252:U:C5	2.56	0.41
1:CA:399:G:H2'	1:CA:400:C:H6	1.84	0.41
1:CA:518:C:OP2	1:CA:530:G:H1'	2.21	0.41
1:CA:58:C:O2'	1:CA:59:A:H5'	2.21	0.41
1:CA:643:C:H2'	1:CA:644:U:C6	2.55	0.41
1:CA:974:A:OP2	21:CN:80:ARG:NH1	2.53	0.41
18:CB:18:GLN:OE1	18:CB:189:ASN:OD1	2.39	0.41
18:CB:56:LEU:H	18:CB:56:LEU:HD12	1.85	0.41
2:CC:66:THR:HG22	2:CC:101:ASN:HD22	1.86	0.41
3:CD:25:ARG:HB2	3:CD:25:ARG:HE	1.71	0.41
6:CG:77:ARG:O	6:CG:84:TYR:N	2.41	0.41
7:CH:71:VAL:HG12	7:CH:72:GLU:N	2.36	0.41
9:CJ:59:LYS:C	9:CJ:61:ALA:H	2.25	0.41
12:CM:1:ALA:H2	12:CM:8:ILE:HG22	1.82	0.41
15:CR:52:ARG:HG3	15:CR:52:ARG:HH11	1.86	0.41
17:CT:53:MET:HA	17:CT:56:ILE:CG1	2.50	0.41
36:D2:3:ARG:HH21	36:D2:3:ARG:HG2	1.86	0.41
23:DB:2742:G:P	32:D4:24:ARG:HH12	2.44	0.41
23:DB:1076:C:H2'	23:DB:1077:A:H8	1.85	0.41
23:DB:1176:U:H3'	23:DB:1177:G:O4'	2.20	0.41
23:DB:1361:G:H2'	23:DB:1362:C:H6	1.85	0.41
23:DB:1373:A:H4'	23:DB:2212:A:N3	2.36	0.41
23:DB:1649:G:O2'	23:DB:1650:A:H5'	2.21	0.41
23:DB:2266:A:H4'	23:DB:2267:A:O5'	2.21	0.41
23:DB:2315:G:H2'	23:DB:2316:G:H8	1.86	0.41
23:DB:389:G:C8	23:DB:2413:G:H4'	2.56	0.41
23:DB:2471:A:O2'	23:DB:2472:G:C5'	2.68	0.41
23:DB:2557:G:H2'	23:DB:2558:C:H6	1.84	0.41
23:DB:2569:G:C2	23:DB:2570:G:C8	3.09	0.41
23:DB:2734:A:H2'	23:DB:2735:G:C5'	2.49	0.41
23:DB:969:G:H2'	23:DB:970:U:H6	1.84	0.41
25:DC:173:LEU:HD22	25:DC:181:ARG:O	2.20	0.41
26:DD:61:THR:HG23	26:DD:64:GLU:OE2	2.21	0.41
29:DE:133:LEU:C	29:DE:135:ALA:H	2.25	0.41
29:DE:118:LEU:HA	29:DE:186:VAL:CG1	2.51	0.41
47:DF:133:GLU:HA	47:DF:150:GLY:HA2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DF:151:LEU:HD12	47:DF:151:LEU:C	2.41	0.41
23:DB:2314:A:H4'	47:DF:34:THR:HG21	2.02	0.41
23:DB:2751:G:OP2	48:DG:2:ARG:HD2	2.21	0.41
24:DI:73:PRO:CG	24:DI:78:LEU:HD21	2.48	0.41
41:DJ:45:THR:N	41:DJ:46:PRO:CD	2.84	0.41
41:DJ:84:ILE:HG13	41:DJ:84:ILE:O	2.21	0.41
38:DM:31:PHE:HB3	38:DM:130:PHE:CE1	2.55	0.41
38:DM:31:PHE:CZ	38:DM:110:GLU:HB3	2.56	0.41
44:DQ:55:GLN:O	44:DQ:59:LEU:N	2.54	0.41
50:DT:31:VAL:C	50:DT:32:LEU:HD23	2.42	0.41
46:DU:50:ALA:H	46:DU:53:GLN:NE2	2.18	0.41
46:DU:73:ASN:HA	46:DU:95:PHE:CZ	2.56	0.41
51:DZ:18:ARG:HA	51:DZ:24:ALA:HA	2.03	0.41
51:DZ:19:SER:OG	51:DZ:23:ASN:HB2	2.20	0.41
1:AA:1237:C:C5'	1:AA:1238:A:H5'	2.51	0.40
1:AA:1239:A:H1'	1:AA:1241:G:C4	2.56	0.40
1:AA:503:C:O2'	1:AA:504:C:H5'	2.20	0.40
1:AA:716:A:H2'	1:AA:717:U:H6	1.85	0.40
1:AA:771:G:H2'	1:AA:772:U:C6	2.56	0.40
1:AA:815:A:H4'	1:AA:817:C:C4	2.56	0.40
1:AA:82:G:C6	1:AA:88:U:O2	2.74	0.40
18:AB:17:HIS:NE2	18:AB:18:GLN:NE2	2.69	0.40
18:AB:184:ALA:O	18:AB:199:ILE:HB	2.21	0.40
18:AB:42:LEU:HA	18:AB:45:THR:CB	2.50	0.40
5:AF:40:GLU:H	5:AF:61:LEU:HB2	1.86	0.40
5:AF:6:ILE:HD11	5:AF:8:PHE:CD2	2.45	0.40
6:AG:60:ALA:O	6:AG:63:VAL:HB	2.21	0.40
1:AA:1385:G:H5'	8:AI:129:ARG:NH2	2.36	0.40
11:AL:35:ARG:HA	11:AL:35:ARG:NH1	2.36	0.40
20:AO:37:ASN:CA	20:AO:40:GLN:HE22	2.33	0.40
14:AQ:44:HIS:HB2	14:AQ:69:THR:O	2.21	0.40
14:AQ:65:PRO:HA	14:AQ:71:SER:OG	2.21	0.40
16:AS:14:LEU:HA	16:AS:17:LYS:NZ	2.36	0.40
17:AT:53:MET:HA	17:AT:56:ILE:HG13	2.03	0.40
23:BB:1159:U:O2'	23:BB:1160:G:H5'	2.22	0.40
23:BB:1228:G:H2'	23:BB:1229:C:H6	1.86	0.40
23:BB:1829:A:H3'	23:BB:1830:C:H6	1.86	0.40
23:BB:195:A:H5''	37:BL:47:ARG:NH2	2.36	0.40
23:BB:2219:U:O2'	23:BB:2220:U:H5'	2.21	0.40
23:BB:2329:U:H2'	23:BB:2330:G:H8	1.85	0.40
23:BB:2371:G:H8	23:BB:2371:G:OP2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2383:G:H5'	23:BB:2383:G:H8	1.85	0.40
23:BB:2064:C:H1'	23:BB:2450:A:C5	2.56	0.40
23:BB:2520:C:O2'	23:BB:2521:C:H5'	2.20	0.40
23:BB:2748:A:O2'	23:BB:2749:A:H5'	2.21	0.40
23:BB:2772:C:H2'	23:BB:2773:C:H6	1.85	0.40
23:BB:527:C:N4	23:BB:2779:U:OP2	2.54	0.40
23:BB:2896:C:H2'	23:BB:2897:U:H6	1.85	0.40
23:BB:321:U:H4'	29:BE:159:LEU:O	2.21	0.40
23:BB:705:A:H8	23:BB:705:A:O5'	2.03	0.40
23:BB:743:A:H2'	23:BB:744:U:C6	2.57	0.40
23:BB:769:U:H2'	23:BB:770:G:C8	2.56	0.40
23:BB:817:C:O2	23:BB:839:U:H4'	2.22	0.40
23:BB:845:A:H2'	23:BB:846:U:C5'	2.39	0.40
25:BC:93:VAL:CG1	25:BC:94:LEU:H	2.19	0.40
29:BE:191:ASP:O	29:BE:194:LYS:HB3	2.21	0.40
48:BG:69:ALA:O	48:BG:73:SER:N	2.52	0.40
40:BH:116:ARG:HB2	40:BH:116:ARG:CZ	2.50	0.40
40:BH:133:GLN:HE21	40:BH:139:PHE:HD2	1.67	0.40
27:BK:111:LYS:HD3	27:BK:111:LYS:N	2.36	0.40
37:BL:81:ASP:O	37:BL:82:LEU:HB2	2.21	0.40
42:BN:37:THR:HB	42:BN:40:LYS:HE2	2.03	0.40
42:BN:82:GLU:C	42:BN:84:GLY:N	2.74	0.40
43:BO:115:LEU:N	43:BO:115:LEU:HD22	2.35	0.40
28:BP:29:VAL:O	28:BP:40:GLN:N	2.51	0.40
44:BQ:91:ARG:HH12	49:BR:10:LYS:HB3	1.86	0.40
49:BR:49:ILE:HG21	49:BR:54:VAL:HA	2.02	0.40
50:BT:69:ARG:HG2	50:BT:74:ILE:CA	2.51	0.40
46:BU:1:ALA:HB1	46:BU:84:PHE:CZ	2.56	0.40
46:BU:18:LYS:O	46:BU:20:LYS:N	2.54	0.40
39:BX:8:GLU:HB3	39:BX:12:GLU:CB	2.51	0.40
30:BY:46:MET:O	30:BY:49:ALA:HB3	2.21	0.40
30:BY:7:THR:CG2	30:BY:8:GLN:N	2.83	0.40
1:CA:1315:U:H3'	1:CA:1316:G:H8	1.85	0.40
1:CA:279:A:H5''	1:CA:280:C:C3'	2.40	0.40
1:CA:684:U:H2'	1:CA:685:G:O4'	2.21	0.40
1:CA:769:G:O2'	1:CA:770:C:H5'	2.21	0.40
1:CA:86:G:H1'	1:CA:87:C:C6	2.56	0.40
2:CC:73:GLY:O	2:CC:77:GLY:N	2.51	0.40
3:CD:110:ARG:HG3	3:CD:110:ARG:NH1	2.35	0.40
4:CE:148:SER:O	4:CE:151:MET:HB2	2.20	0.40
6:CG:75:LYS:HA	6:CG:75:LYS:HZ2	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:7:ARG:NE	9:CJ:101:SER:HB3	2.34	0.40
11:CL:23:LEU:C	11:CL:25:ALA:N	2.75	0.40
1:CA:1047:G:O3'	21:CN:3:GLN:HB3	2.21	0.40
16:CS:14:LEU:HD12	16:CS:15:LEU:N	2.36	0.40
23:DB:1098:A:O3'	24:DI:4:VAL:O	2.39	0.40
23:DB:1135:C:N4	23:DB:1139:G:C6	2.89	0.40
23:DB:1279:G:H2'	23:DB:1280:G:O4'	2.21	0.40
23:DB:123:G:H4'	23:DB:1376:C:O5'	2.21	0.40
23:DB:1433:A:H2'	23:DB:1434:A:O4'	2.21	0.40
23:DB:146:A:H2'	23:DB:147:C:H6	1.84	0.40
23:DB:14:A:N6	23:DB:15:G:C2	2.90	0.40
23:DB:2247:A:H3'	56:DB:3575:HOH:O	2.21	0.40
23:DB:295:G:O2'	23:DB:296:U:H5'	2.21	0.40
23:DB:357:C:H2'	23:DB:358:U:H6	1.86	0.40
23:DB:454:A:H3'	23:DB:455:C:C5'	2.47	0.40
23:DB:523:C:O2'	23:DB:524:G:H5'	2.20	0.40
23:DB:812:C:H2'	23:DB:813:U:H6	1.85	0.40
26:DD:202:ILE:HG22	26:DD:204:LYS:HZ1	1.86	0.40
48:DG:116:LEU:CD2	48:DG:122:ALA:H	2.34	0.40
48:DG:132:LEU:CD2	48:DG:132:LEU:H	2.27	0.40
24:DI:10:LEU:C	24:DI:10:LEU:HD12	2.41	0.40
41:DJ:72:LYS:O	41:DJ:73:VAL:HG13	2.21	0.40
38:DM:31:PHE:O	38:DM:104:GLU:HA	2.21	0.40
38:DM:66:ARG:HB2	38:DM:101:VAL:HG13	2.04	0.40
43:DO:89:ASP:H	43:DO:115:LEU:HA	1.86	0.40
49:DR:82:HIS:O	49:DR:83:TYR:C	2.58	0.40
50:DT:40:LYS:CA	50:DT:43:ILE:HG22	2.51	0.40
35:DV:2:PHE:CE2	35:DV:55:GLU:HB3	2.56	0.40
52:DW:75:ASN:O	52:DW:76:ARG:HB2	2.21	0.40
23:DB:77:G:H5'	39:DX:52:ARG:HG2	2.03	0.40
1:AA:17:U:H4'	1:AA:1080:A:O4'	2.20	0.40
1:AA:1303:C:O2'	1:AA:1304:G:H5'	2.21	0.40
1:AA:1322:C:O4'	1:AA:1322:C:O2	2.36	0.40
1:AA:1426:G:H2'	1:AA:1427:C:O4'	2.21	0.40
1:AA:279:A:H5''	1:AA:280:C:C3'	2.44	0.40
1:AA:59:A:H1'	1:AA:354:G:N2	2.35	0.40
1:AA:366:A:H1'	1:AA:395:C:O2	2.21	0.40
1:AA:659:U:O2'	1:AA:660:C:H5'	2.21	0.40
1:AA:715:A:H8	1:AA:715:A:O5'	2.05	0.40
18:AB:60:ALA:O	18:AB:224:ARG:HB3	2.21	0.40
18:AB:73:ARG:CZ	18:AB:73:ARG:HB2	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:168:ARG:HE	2:AC:169:GLU:CA	2.34	0.40
2:AC:168:ARG:HG2	2:AC:169:GLU:H	1.86	0.40
2:AC:60:ALA:C	2:AC:62:SER:H	2.23	0.40
4:AE:106:ALA:HB1	4:AE:110:MET:CB	2.50	0.40
5:AF:4:TYR:HA	5:AF:90:MET:O	2.21	0.40
6:AG:145:GLU:CA	6:AG:148:LYS:HB2	2.42	0.40
7:AH:10:LEU:O	7:AH:13:ILE:HB	2.21	0.40
8:AI:117:LEU:HD13	8:AI:121:ARG:O	2.21	0.40
8:AI:17:ARG:NH2	8:AI:65:THR:HG21	2.35	0.40
12:AM:19:THR:HA	12:AM:24:VAL:CG2	2.48	0.40
21:AN:32:ASP:OD2	21:AN:33:VAL:N	2.54	0.40
21:AN:80:ARG:HG2	21:AN:80:ARG:HH11	1.86	0.40
19:AU:40:PRO:O	19:AU:44:ARG:HB2	2.21	0.40
33:B1:34:GLU:CD	33:B1:49:LYS:HG3	2.41	0.40
36:B2:22:MET:HG2	36:B2:31:LEU:CD1	2.51	0.40
34:B3:44:ARG:N	34:B3:45:PRO:CD	2.81	0.40
22:BA:30:C:H2'	22:BA:30:C:O2	2.20	0.40
22:BA:63:C:H2'	22:BA:64:G:H8	1.87	0.40
23:BB:1019:U:H2'	23:BB:1020:A:H8	1.77	0.40
23:BB:1128:G:C6	23:BB:2518:A:N6	2.89	0.40
23:BB:1381:G:H1'	23:BB:1571:A:N1	2.36	0.40
23:BB:1463:C:H2'	23:BB:1464:G:C8	2.56	0.40
23:BB:1473:G:O2'	23:BB:1474:U:H5'	2.21	0.40
23:BB:1937:A:C8	23:BB:1939:U:H2'	2.55	0.40
23:BB:1997:C:O2'	23:BB:1998:A:H5'	2.20	0.40
23:BB:2155:U:H2'	23:BB:2156:G:C8	2.56	0.40
23:BB:2347:C:N3	23:BB:2371:G:N2	2.70	0.40
23:BB:2450:A:O2'	23:BB:2451:A:H5'	2.21	0.40
23:BB:2778:A:H4'	23:BB:2779:U:OP2	2.21	0.40
23:BB:496:G:H4'	45:BS:61:ASN:OD1	2.20	0.40
23:BB:662:G:O2'	23:BB:663:G:H5'	2.20	0.40
23:BB:848:C:H2'	23:BB:849:A:H8	1.83	0.40
23:BB:853:C:H2'	23:BB:854:C:H6	1.86	0.40
23:BB:970:U:O2'	23:BB:984:A:H4'	2.21	0.40
25:BC:171:VAL:HG23	25:BC:185:ALA:CB	2.52	0.40
25:BC:93:VAL:HG12	25:BC:101:ARG:O	2.22	0.40
26:BD:182:ALA:C	26:BD:184:ARG:N	2.75	0.40
29:BE:24:ASN:OD1	29:BE:27:LEU:HB2	2.21	0.40
29:BE:37:ALA:O	29:BE:39:ALA:N	2.47	0.40
47:BF:100:GLU:C	47:BF:102:LEU:N	2.74	0.40
47:BF:148:VAL:HG23	47:BF:149:ARG:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BF:51:ASN:C	47:BF:53:ALA:N	2.74	0.40
47:BF:68:LYS:N	47:BF:68:LYS:HD2	2.36	0.40
47:BF:86:CYS:O	47:BF:88:VAL:HG23	2.22	0.40
48:BG:94:ARG:NH2	48:BG:104:LEU:HA	2.34	0.40
48:BG:153:PRO:HB3	48:BG:158:GLY:CA	2.46	0.40
40:BH:83:LYS:HZ2	40:BH:83:LYS:HA	1.87	0.40
41:BJ:29:ALA:O	41:BJ:32:LEU:HB2	2.21	0.40
41:BJ:81:ILE:HG23	41:BJ:82:GLY:N	2.25	0.40
27:BK:62:VAL:HG12	27:BK:63:VAL:H	1.86	0.40
27:BK:67:LYS:C	27:BK:67:LYS:HD3	2.41	0.40
37:BL:14:LYS:HB3	37:BL:14:LYS:HE3	1.96	0.40
37:BL:23:ILE:O	37:BL:24:GLY:C	2.60	0.40
38:BM:63:ILE:N	38:BM:63:ILE:HD12	2.36	0.40
28:BP:89:GLY:N	28:BP:112:ARG:NH1	2.68	0.40
49:BR:7:SER:HB2	49:BR:22:LEU:HD13	2.03	0.40
45:BS:21:ALA:HB1	45:BS:74:ILE:HG12	2.03	0.40
35:BV:35:GLU:HB2	35:BV:93:ARG:HD2	2.03	0.40
51:BZ:54:LYS:O	51:BZ:57:ARG:HB2	2.20	0.40
1:CA:1011:C:H2'	1:CA:1012:A:H8	1.87	0.40
1:CA:1118:U:N3	1:CA:1156:G:N2	2.69	0.40
1:CA:123:U:H2'	1:CA:124:C:C6	2.56	0.40
1:CA:1240:U:H3'	1:CA:1241:G:H5'	2.03	0.40
1:CA:345:C:H3'	28:DP:38:ARG:NH1	2.36	0.40
1:CA:373:A:H2'	1:CA:374:A:C8	2.52	0.40
1:CA:505:G:C6	1:CA:535:A:C2	3.10	0.40
1:CA:647:C:H2'	1:CA:648:A:C8	2.57	0.40
1:CA:667:G:H2'	1:CA:668:G:C8	2.55	0.40
1:CA:909:A:H2	1:CA:1413:A:N3	2.20	0.40
18:CB:218:ALA:HA	18:CB:221:ARG:NH2	2.35	0.40
18:CB:58:LYS:HD3	18:CB:59:ILE:HG23	2.03	0.40
18:CB:64:GLY:C	18:CB:87:ASP:O	2.60	0.40
2:CC:102:ILE:HD12	2:CC:102:ILE:N	2.37	0.40
2:CC:111:ASP:HA	2:CC:182:ASP:OD1	2.21	0.40
2:CC:72:PRO:C	2:CC:76:ILE:HG12	2.41	0.40
3:CD:144:ILE:HG22	3:CD:145:ARG:H	1.86	0.40
3:CD:171:GLU:HG3	3:CD:182:LYS:HD2	2.03	0.40
4:CE:152:VAL:CG1	4:CE:156:ARG:HE	2.35	0.40
7:CH:17:GLN:HE21	7:CH:62:LEU:HD23	1.86	0.40
8:CI:108:ARG:HG2	8:CI:108:ARG:H	1.59	0.40
8:CI:117:LEU:HB3	8:CI:122:ARG:O	2.21	0.40
9:CJ:7:ARG:HH21	9:CJ:101:SER:CB	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:CL:115:LYS:O	11:CL:116:TYR:HB2	2.21	0.40
21:CN:80:ARG:HG2	21:CN:81:ILE:N	2.36	0.40
33:D1:32:LYS:O	33:D1:34:GLU:HG3	2.21	0.40
36:D2:22:MET:HG2	36:D2:31:LEU:CD1	2.51	0.40
34:D3:50:SER:C	34:D3:52:GLY:N	2.74	0.40
23:DB:1286:A:C6	23:DB:1329:U:C2	3.09	0.40
23:DB:1567:G:H5'	25:DC:84:PRO:HB3	2.04	0.40
23:DB:1848:A:H2'	23:DB:1849:G:C8	2.56	0.40
23:DB:2679:A:H2'	23:DB:2680:U:C6	2.57	0.40
23:DB:2850:A:O4'	23:DB:2868:A:H2	2.04	0.40
23:DB:296:U:H2'	23:DB:297:G:H8	1.87	0.40
23:DB:724:U:O2'	23:DB:725:G:H5'	2.21	0.40
23:DB:732:C:H2'	23:DB:733:G:O4'	2.21	0.40
23:DB:780:G:H2'	23:DB:782:A:N7	2.36	0.40
23:DB:853:C:H2'	23:DB:854:C:H6	1.87	0.40
25:DC:100:ARG:H	25:DC:100:ARG:HG2	1.75	0.40
25:DC:29:PHE:CE2	25:DC:31:PRO:HG2	2.56	0.40
26:DD:8:LYS:HG2	26:DD:197:THR:N	2.36	0.40
29:DE:12:LEU:CD1	29:DE:14:VAL:HG13	2.51	0.40
29:DE:138:LEU:O	29:DE:142:ALA:N	2.54	0.40
29:DE:191:ASP:O	29:DE:194:LYS:HB3	2.22	0.40
29:DE:61:ARG:HB3	29:DE:61:ARG:CZ	2.51	0.40
47:DF:104:THR:C	47:DF:108:PRO:HG2	2.41	0.40
48:DG:6:ALA:HA	48:DG:7:PRO:HD3	1.82	0.40
48:DG:93:TYR:O	48:DG:94:ARG:HG3	2.21	0.40
40:DH:6:LEU:O	40:DH:15:LEU:HD22	2.21	0.40
40:DH:2:GLN:O	40:DH:19:VAL:O	2.39	0.40
40:DH:50:ARG:C	40:DH:52:ALA:N	2.74	0.40
24:DI:63:ASP:O	24:DI:63:ASP:OD1	2.39	0.40
41:DJ:77:HIS:ND1	41:DJ:79:GLY:N	2.69	0.40
27:DK:41:ILE:CG1	27:DK:42:THR:N	2.84	0.40
28:DP:99:LEU:HA	28:DP:102:ARG:HG3	2.02	0.40
28:DP:53:GLY:O	28:DP:54:LEU:C	2.60	0.40
28:DP:19:PHE:CE2	28:DP:83:ILE:HD11	2.55	0.40
45:DS:54:ALA:C	45:DS:56:ALA:N	2.75	0.40
35:DV:38:LEU:HD12	35:DV:39:ALA:N	2.36	0.40
35:DV:68:LYS:N	35:DV:68:LYS:HD3	2.34	0.40
39:DX:10:SER:O	39:DX:11:VAL:C	2.60	0.40
39:DX:6:LEU:HD22	39:DX:6:LEU:H	1.86	0.40
1:AA:1418:A:H61	1:AA:1482:G:H1'	1.85	0.40
1:AA:230:G:O2'	1:AA:231:U:H5'	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:238:A:H2'	1:AA:239:U:C5'	2.48	0.40
1:AA:277:C:H2'	1:AA:278:G:H8	1.86	0.40
1:AA:373:A:H2'	1:AA:374:A:C8	2.53	0.40
2:AC:179:ALA:O	2:AC:180:ASP:C	2.60	0.40
3:AD:24:VAL:CG1	3:AD:160:LEU:HB3	2.51	0.40
5:AF:53:LYS:N	5:AF:53:LYS:HZ3	2.19	0.40
6:AG:129:ASN:C	6:AG:134:VAL:HG21	2.42	0.40
8:AI:113:LYS:HA	8:AI:120:ALA:CA	2.51	0.40
8:AI:56:MET:C	8:AI:58:GLU:N	2.75	0.40
8:AI:20:ILE:HG23	8:AI:60:LEU:HD12	2.04	0.40
11:AL:115:LYS:O	11:AL:116:TYR:HB2	2.21	0.40
12:AM:42:VAL:HB	12:AM:47:LEU:HD22	2.03	0.40
12:AM:61:LYS:N	12:AM:61:LYS:HE2	2.36	0.40
23:BB:2348:U:H1'	33:B1:38:PHE:HE2	1.86	0.40
33:B1:26:LYS:HG2	33:B1:52:LYS:HD2	2.04	0.40
23:BB:1029:A:H8	23:BB:1029:A:O5'	2.04	0.40
23:BB:1228:G:O2'	23:BB:1229:C:H5'	2.22	0.40
23:BB:1344:U:H5'	23:BB:1384:A:C6	2.57	0.40
23:BB:1729:U:C5	23:BB:1731:G:N2	2.88	0.40
23:BB:1827:U:C2'	23:BB:1828:G:H5'	2.52	0.40
23:BB:2410:G:H2'	23:BB:2411:A:C8	2.56	0.40
23:BB:2032:G:N7	23:BB:2454:G:H1'	2.36	0.40
23:BB:2749:A:OP1	23:BB:2751:G:H5'	2.22	0.40
23:BB:786:C:H5''	23:BB:1780:A:N7	2.36	0.40
23:BB:834:G:O2'	23:BB:835:C:H5'	2.20	0.40
23:BB:860:U:O2'	23:BB:861:A:H5'	2.21	0.40
23:BB:866:A:H61	23:BB:913:U:C1'	2.34	0.40
25:BC:38:LYS:HD2	25:BC:38:LYS:HA	1.95	0.40
25:BC:75:ALA:HB1	25:BC:93:VAL:HG22	2.03	0.40
47:BF:121:PHE:N	47:BF:121:PHE:CD2	2.89	0.40
47:BF:34:THR:OG1	47:BF:154:THR:N	2.53	0.40
47:BF:36:ASN:O	47:BF:86:CYS:O	2.38	0.40
48:BG:34:ARG:HG2	48:BG:34:ARG:HH11	1.85	0.40
40:BH:96:THR:HB	40:BH:112:LYS:O	2.21	0.40
40:BH:2:GLN:O	40:BH:19:VAL:O	2.40	0.40
24:BI:27:LEU:HB2	24:BI:32:VAL:HG21	2.02	0.40
24:BI:49:GLU:OE1	24:BI:52:LEU:HD22	2.21	0.40
41:BJ:26:GLY:C	41:BJ:28:LEU:N	2.75	0.40
27:BK:8:LEU:N	27:BK:8:LEU:HD12	2.36	0.40
43:BO:66:GLY:CA	43:BO:102:ARG:HH21	2.19	0.40
41:BJ:44:TYR:HB2	44:BQ:63:ARG:CD	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BS:49:LYS:HA	45:BS:52:GLU:HG2	2.02	0.40
23:BB:485:C:O2'	45:BS:60:HIS:NE2	2.53	0.40
45:BS:74:ILE:O	45:BS:75:PHE:HB3	2.22	0.40
45:BS:9:HIS:H	45:BS:102:HIS:CE1	2.38	0.40
50:BT:31:VAL:O	50:BT:32:LEU:HB3	2.20	0.40
50:BT:74:ILE:HG13	50:BT:75:GLY:H	1.86	0.40
50:BT:9:LYS:HZ3	50:BT:9:LYS:HB3	1.84	0.40
35:BV:24:ASN:HB3	35:BV:44:HIS:HB3	2.03	0.40
35:BV:60:VAL:HG12	35:BV:61:LEU:N	2.36	0.40
23:BB:851:C:O4'	30:BY:46:MET:HG2	2.21	0.40
51:BZ:56:MET:C	51:BZ:58:VAL:N	2.72	0.40
1:CA:1096:C:H2'	1:CA:1097:C:H6	1.87	0.40
1:CA:1127:G:O2'	1:CA:1128:C:H5'	2.21	0.40
1:CA:1179:A:H2'	1:CA:1180:A:H8	1.87	0.40
1:CA:1238:A:N3	1:CA:1238:A:H2'	2.36	0.40
1:CA:1453:G:H3'	1:CA:1453:G:N3	2.37	0.40
1:CA:123:U:OP1	1:CA:312:C:H5'	2.21	0.40
2:CC:33:ASP:O	2:CC:36:PHE:HB3	2.21	0.40
1:CA:620:C:C2	3:CD:131:ILE:HG21	2.57	0.40
3:CD:143:SER:OG	3:CD:144:ILE:N	2.54	0.40
3:CD:152:SER:CA	3:CD:155:LYS:HD3	2.36	0.40
4:CE:9:GLU:O	4:CE:40:ASP:HA	2.22	0.40
7:CH:68:LYS:HG3	7:CH:69:ALA:O	2.21	0.40
8:CI:44:ARG:O	8:CI:48:ARG:HG3	2.21	0.40
9:CJ:57:VAL:HG22	9:CJ:58:ASN:N	2.28	0.40
9:CJ:8:ILE:O	9:CJ:8:ILE:HG22	2.21	0.40
11:CL:49:ARG:CG	11:CL:89:LEU:HD21	2.49	0.40
21:CN:70:HIS:O	21:CN:71:GLY:C	2.58	0.40
9:CJ:65:TYR:C	21:CN:98:ALA:HB2	2.41	0.40
20:CO:47:LYS:O	20:CO:49:ASP:N	2.54	0.40
13:CP:22:ALA:CB	13:CP:32:PHE:HA	2.51	0.40
13:CP:70:ARG:O	13:CP:73:ALA:HB3	2.22	0.40
16:CS:38:THR:CG2	16:CS:39:ILE:N	2.85	0.40
16:CS:39:ILE:HG21	16:CS:65:MET:O	2.21	0.40
32:D4:15:LYS:O	32:D4:16:ILE:CB	2.69	0.40
23:DB:1146:C:H2'	23:DB:1147:A:C8	2.57	0.40
23:DB:1197:G:H5'	23:DB:1227:G:O2'	2.22	0.40
23:DB:1346:G:C6	23:DB:1601:G:C6	3.09	0.40
23:DB:1523:U:H5''	23:DB:1524:G:H8	1.87	0.40
23:DB:1657:U:C2'	23:DB:1658:C:H5'	2.51	0.40
23:DB:1678:A:H2'	23:DB:1679:A:H8	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2109:U:O2'	23:DB:2110:G:H5'	2.21	0.40
23:DB:245:G:H2'	23:DB:246:C:H6	1.86	0.40
23:DB:2472:G:C8	23:DB:2475:C:N4	2.89	0.40
23:DB:2560:A:H2'	23:DB:2561:U:C6	2.56	0.40
23:DB:2699:C:H2'	23:DB:2700:A:C8	2.56	0.40
23:DB:2746:U:H5''	48:DG:137:LYS:HB3	2.04	0.40
23:DB:2800:A:N3	23:DB:2801:G:H1'	2.37	0.40
23:DB:2822:G:O5'	23:DB:2822:G:H8	2.05	0.40
23:DB:296:U:H2'	23:DB:297:G:C8	2.57	0.40
23:DB:476:G:H4'	23:DB:502:A:H61	1.86	0.40
23:DB:621:A:H2'	23:DB:622:G:O4'	2.22	0.40
23:DB:636:G:H5'	23:DB:639:U:OP1	2.22	0.40
25:DC:36:ASN:HD22	25:DC:59:GLN:C	2.24	0.40
26:DD:141:ARG:O	26:DD:141:ARG:HG3	2.21	0.40
26:DD:56:LYS:C	26:DD:58:ASN:N	2.74	0.40
23:DB:674:G:O3'	29:DE:60:TRP:HZ2	2.04	0.40
47:DF:135:ILE:C	47:DF:137:PHE:N	2.75	0.40
40:DH:28:ASN:HA	40:DH:28:ASN:HD22	1.61	0.40
41:DJ:28:LEU:HD23	41:DJ:29:ALA:H	1.86	0.40
27:DK:47:ILE:CG1	27:DK:48:PRO:HD2	2.36	0.40
27:DK:53:LYS:HB2	27:DK:54:LYS:HD2	2.03	0.40
37:DL:89:VAL:O	37:DL:91:ASP:N	2.55	0.40
38:DM:10:ARG:HG3	38:DM:10:ARG:NH2	2.36	0.40
38:DM:111:GLU:HA	38:DM:114:ARG:NH2	2.35	0.40
38:DM:116:ALA:HA	38:DM:119:LEU:HD12	2.03	0.40
23:DB:2723:C:H5''	42:DN:1:MET:HE2	2.03	0.40
42:DN:82:GLU:C	42:DN:84:GLY:H	2.25	0.40
28:DP:74:GLN:O	28:DP:75:THR:C	2.59	0.40
50:DT:31:VAL:O	50:DT:32:LEU:HB3	2.21	0.40
52:DW:70:VAL:O	52:DW:70:VAL:HG22	2.21	0.40
39:DX:2:LYS:O	39:DX:5:GLU:HG2	2.21	0.40
1:AA:1041:G:H2'	1:AA:1042:A:O4'	2.22	0.40
1:AA:35:G:H2'	1:AA:36:C:C6	2.56	0.40
1:AA:57:G:N2	1:AA:388:G:C6	2.89	0.40
1:AA:419:C:O2'	1:AA:420:U:H5'	2.21	0.40
1:AA:588:G:H2'	1:AA:588:G:N3	2.36	0.40
1:AA:570:G:H1'	1:AA:820:U:C4	2.56	0.40
1:AA:923:A:O5'	1:AA:923:A:H8	2.05	0.40
1:AA:931:C:H2'	1:AA:932:C:C6	2.57	0.40
1:AA:957:U:H3	1:AA:960:U:H5''	1.87	0.40
1:AA:957:U:N3	1:AA:960:U:OP2	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AB:45:THR:HG23	18:AB:200:PRO:O	2.21	0.40
2:AC:109:GLU:O	2:AC:140:ALA:HA	2.20	0.40
2:AC:137:VAL:HA	2:AC:148:ILE:CD1	2.51	0.40
2:AC:142:ARG:C	2:AC:144:GLY:H	2.24	0.40
2:AC:194:VAL:CG1	2:AC:195:ILE:N	2.84	0.40
2:AC:182:ASP:O	2:AC:201:ILE:N	2.54	0.40
2:AC:48:LYS:CD	2:AC:48:LYS:H	2.34	0.40
2:AC:51:VAL:HG22	2:AC:52:SER:N	2.37	0.40
3:AD:24:VAL:HG23	3:AD:25:ARG:N	2.36	0.40
4:AE:37:VAL:HG12	4:AE:47:PHE:CB	2.51	0.40
5:AF:85:ILE:HG22	5:AF:86:ARG:H	1.87	0.40
1:AA:1349:A:OP1	8:AI:121:ARG:HB2	2.22	0.40
8:AI:12:LYS:C	8:AI:14:SER:H	2.25	0.40
8:AI:5:TYR:O	8:AI:19:PHE:HA	2.21	0.40
9:AJ:59:LYS:C	9:AJ:61:ALA:H	2.24	0.40
12:AM:96:VAL:C	12:AM:98:GLY:H	2.24	0.40
21:AN:5:MET:HE3	21:AN:60:ARG:NH1	2.37	0.40
21:AN:92:ILE:HG21	21:AN:95:LEU:CD2	2.52	0.40
13:AP:39:PHE:CG	13:AP:40:ASN:N	2.89	0.40
13:AP:71:VAL:HA	13:AP:74:LEU:HG	2.02	0.40
17:AT:24:ARG:HG3	17:AT:65:LEU:HD21	2.04	0.40
32:B4:15:LYS:O	32:B4:16:ILE:CB	2.69	0.40
32:B4:25:VAL:O	32:B4:26:ILE:HD13	2.21	0.40
22:BA:65:U:C2'	22:BA:66:A:H5'	2.52	0.40
23:BB:1068:G:C6	23:BB:1069:A:N6	2.89	0.40
23:BB:1301:A:N3	23:BB:1301:A:H2'	2.35	0.40
23:BB:1759:A:H2'	23:BB:1759:A:N3	2.37	0.40
23:BB:1942:C:H6	23:BB:1942:C:O5'	2.04	0.40
23:BB:2025:C:H2'	23:BB:2026:U:C6	2.57	0.40
23:BB:2389:G:H5''	23:BB:2390:U:O4'	2.21	0.40
23:BB:2419:U:H2'	23:BB:2420:C:C6	2.57	0.40
23:BB:2651:C:H2'	23:BB:2652:C:H6	1.87	0.40
23:BB:2821:A:OP2	23:BB:2822:G:OP2	2.40	0.40
23:BB:479:A:O2'	23:BB:481:G:H5'	2.20	0.40
23:BB:744:U:H2'	23:BB:745:G:C8	2.56	0.40
23:BB:856:G:H2'	23:BB:857:G:C8	2.57	0.40
25:BC:129:LEU:HB3	25:BC:134:ILE:CG2	2.52	0.40
25:BC:69:ASN:O	25:BC:70:LYS:C	2.59	0.40
26:BD:102:ALA:C	26:BD:104:VAL:H	2.24	0.40
26:BD:191:GLY:O	26:BD:192:ALA:HB3	2.22	0.40
26:BD:90:PHE:N	26:BD:94:GLN:OE1	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BE:17:THR:C	29:BE:19:PHE:N	2.74	0.40
29:BE:61:ARG:HB3	29:BE:61:ARG:CZ	2.52	0.40
48:BG:123:GLU:C	48:BG:125:PRO:HD3	2.42	0.40
48:BG:29:ASN:HD22	48:BG:30:GLY:N	2.19	0.40
48:BG:88:LEU:HD12	48:BG:88:LEU:C	2.42	0.40
42:BN:114:GLU:HG2	42:BN:115:LEU:O	2.21	0.40
42:BN:60:VAL:O	42:BN:60:VAL:HG12	2.21	0.40
44:BQ:48:ASP:OD2	44:BQ:48:ASP:N	2.53	0.40
52:BW:77:LYS:HD3	52:BW:77:LYS:HA	1.84	0.40
51:BZ:5:CYS:CB	51:BZ:10:LYS:N	2.85	0.40
51:BZ:21:ALA:HB3	51:BZ:23:ASN:HD21	1.85	0.40
1:CA:1011:C:H2'	1:CA:1012:A:C8	2.56	0.40
1:CA:1085:U:H3'	1:CA:1086:U:C5	2.56	0.40
1:CA:1096:C:H2'	1:CA:1097:C:C6	2.56	0.40
1:CA:1098:C:H2'	1:CA:1099:G:O4'	2.22	0.40
1:CA:1128:C:H4'	1:CA:1148:U:C2	2.57	0.40
1:CA:1176:A:H2'	1:CA:1177:G:C8	2.56	0.40
1:CA:1332:A:H2'	1:CA:1333:A:O4'	2.21	0.40
1:CA:143:A:O3'	1:CA:144:G:H8	2.04	0.40
1:CA:1518:A:H2'	1:CA:1519:A:C8	2.57	0.40
1:CA:1533:C:H2'	1:CA:1533:C:O2	2.22	0.40
1:CA:479:U:O2'	1:CA:480:U:H5'	2.22	0.40
1:CA:592:G:O2'	1:CA:593:U:H5'	2.21	0.40
1:CA:723:U:H3'	1:CA:724:G:C5'	2.52	0.40
1:CA:808:C:O2'	1:CA:809:G:H5'	2.21	0.40
1:CA:844:G:H2'	1:CA:845:A:C8	2.47	0.40
1:CA:928:G:H2'	1:CA:929:G:C8	2.56	0.40
2:CC:148:ILE:HA	2:CC:200:TRP:O	2.22	0.40
2:CC:133:MET:CE	2:CC:152:VAL:HG23	2.51	0.40
3:CD:139:ASN:N	3:CD:181:PHE:O	2.53	0.40
4:CE:132:PRO:O	4:CE:133:ILE:C	2.60	0.40
5:CF:29:ILE:HG21	5:CF:64:VAL:CG1	2.49	0.40
5:CF:3:HIS:CE1	5:CF:94:HIS:HA	2.56	0.40
6:CG:138:GLU:O	6:CG:139:ASP:C	2.60	0.40
6:CG:15:PRO:HG2	6:CG:43:TYR:OH	2.21	0.40
10:CK:60:PHE:CZ	6:CG:146:ALA:HB1	2.56	0.40
12:CM:12:LYS:HD3	12:CM:12:LYS:HA	1.90	0.40
21:CN:66:THR:HG23	21:CN:82:LYS:NZ	2.37	0.40
16:CS:51:HIS:CA	16:CS:56:HIS:HA	2.39	0.40
17:CT:2:ASN:HD22	17:CT:3:ILE:HG13	1.86	0.40
17:CT:50:PHE:O	17:CT:53:MET:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:D3:7:ARG:HA	34:D3:7:ARG:HD2	1.89	0.40
32:D4:30:GLU:HA	32:D4:31:PRO:HD3	1.89	0.40
22:DA:13:G:N2	22:DA:16:G:N3	2.69	0.40
22:DA:65:U:C2'	22:DA:66:A:H5'	2.52	0.40
23:DB:1061:U:O4	24:DI:11:GLN:HG3	2.22	0.40
23:DB:1435:G:H2'	23:DB:1436:G:C8	2.56	0.40
23:DB:1794:A:O2'	23:DB:1795:C:H5'	2.22	0.40
23:DB:1869:G:N2	23:DB:1871:A:C8	2.90	0.40
23:DB:1956:U:O2'	23:DB:1957:C:H5'	2.22	0.40
23:DB:2285:C:H4'	23:DB:2288:A:N6	2.36	0.40
23:DB:2294:G:P	43:DO:94:ARG:NH1	2.94	0.40
23:DB:2320:U:O2'	23:DB:2322:A:N7	2.48	0.40
23:DB:2330:G:N3	52:DW:38:ARG:HB3	2.37	0.40
23:DB:2469:A:C2	23:DB:2482:A:H1'	2.57	0.40
23:DB:379:G:N1	23:DB:396:G:C6	2.90	0.40
23:DB:441:U:O2'	23:DB:442:G:H5'	2.22	0.40
23:DB:458:G:O2'	23:DB:459:U:P	2.79	0.40
23:DB:635:C:H2'	23:DB:636:G:H8	1.86	0.40
23:DB:727:A:O2'	23:DB:728:G:H5'	2.21	0.40
23:DB:817:C:O2'	23:DB:839:U:H5''	2.22	0.40
25:DC:142:ASN:CG	25:DC:142:ASN:O	2.59	0.40
25:DC:4:LYS:CG	25:DC:5:CYS:H	2.32	0.40
29:DE:133:LEU:C	29:DE:135:ALA:N	2.75	0.40
29:DE:153:LEU:HD12	29:DE:154:ASP:N	2.37	0.40
47:DF:131:VAL:C	47:DF:133:GLU:N	2.73	0.40
40:DH:113:SER:HB3	40:DH:114:GLU:H	1.66	0.40
40:DH:31:VAL:O	40:DH:32:PRO:C	2.58	0.40
24:DI:90:GLY:C	24:DI:91:LYS:HD2	2.41	0.40
37:DL:111:ILE:CG2	37:DL:112:LEU:H	2.28	0.40
37:DL:68:SER:C	37:DL:70:LYS:N	2.74	0.40
42:DN:87:PHE:HZ	42:DN:115:LEU:HB3	1.87	0.40
42:DN:119:SER:O	42:DN:120:GLU:HB2	2.21	0.40
42:DN:59:SER:C	42:DN:61:ALA:H	2.25	0.40
43:DO:14:ALA:C	43:DO:16:ARG:N	2.74	0.40
28:DP:85:VAL:O	28:DP:87:ARG:N	2.55	0.40
44:DQ:49:ARG:O	44:DQ:53:LYS:HE2	2.21	0.40
44:DQ:59:LEU:O	44:DQ:59:LEU:HD22	2.21	0.40
44:DQ:85:ALA:O	44:DQ:86:SER:C	2.60	0.40
49:DR:47:VAL:HG12	49:DR:47:VAL:O	2.22	0.40
46:DU:33:VAL:O	46:DU:63:ALA:HB1	2.22	0.40
35:DV:51:GLN:HA	35:DV:56:PHE:CD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DX:17:GLU:O	39:DX:20:ASN:HB2	2.22	0.40
39:DX:28:LEU:HD22	39:DX:37:LEU:HD11	2.02	0.40
1:AA:1171:A:O2'	1:AA:1172:C:H5'	2.21	0.40
1:AA:1346:A:N7	6:AG:9:ARG:NH2	2.70	0.40
1:AA:1479:C:H2'	1:AA:1480:A:H8	1.87	0.40
1:AA:20:U:O2'	1:AA:21:G:H5'	2.21	0.40
1:AA:341:C:O2'	1:AA:342:C:H5'	2.22	0.40
1:AA:628:G:O2'	1:AA:629:A:H5'	2.21	0.40
4:AE:148:SER:HB2	4:AE:150:GLU:OE2	2.21	0.40
4:AE:152:VAL:CG1	4:AE:156:ARG:HE	2.33	0.40
4:AE:56:PRO:O	4:AE:57:ALA:C	2.60	0.40
4:AE:61:LYS:O	4:AE:65:LYS:HG2	2.21	0.40
4:AE:46:GLY:HA3	4:AE:70:MET:HG2	2.03	0.40
5:AF:81:ASN:O	5:AF:83:ALA:N	2.55	0.40
8:AI:20:ILE:HG13	8:AI:62:LEU:CD1	2.48	0.40
12:AM:13:HIS:H	12:AM:16:ILE:CG2	2.34	0.40
15:AR:19:GLU:HG3	15:AR:54:LEU:HD12	2.02	0.40
15:AR:55:ALA:O	15:AR:58:ILE:HB	2.21	0.40
16:AS:48:ILE:HG22	16:AS:49:ALA:N	2.36	0.40
31:B0:50:GLY:O	31:B0:51:ARG:C	2.60	0.40
32:B4:16:ILE:CG1	32:B4:25:VAL:HG22	2.51	0.40
23:BB:104:A:H2'	23:BB:105:C:C6	2.56	0.40
23:BB:1050:A:C2	23:BB:1051:G:H1'	2.56	0.40
23:BB:1062:G:H2'	23:BB:1063:G:H8	1.85	0.40
23:BB:1198:U:O2'	44:BQ:3:VAL:HG13	2.22	0.40
23:BB:1251:C:OP2	44:BQ:9:ALA:HB2	2.22	0.40
23:BB:1341:G:H3'	23:BB:1397:U:O2	2.22	0.40
23:BB:1589:U:H2'	23:BB:1590:A:H8	1.85	0.40
23:BB:1592:C:H2'	23:BB:1593:A:H8	1.87	0.40
23:BB:1900:A:N1	23:BB:1970:A:C6	2.90	0.40
23:BB:192:C:H2'	23:BB:193:U:H5'	2.04	0.40
23:BB:1987:A:H2'	23:BB:1988:G:C8	2.56	0.40
23:BB:2357:G:H22	23:BB:2359:C:H3'	1.86	0.40
23:BB:2773:C:H5''	26:BD:169:ARG:CB	2.49	0.40
23:BB:2850:A:O4'	23:BB:2868:A:H2	2.05	0.40
23:BB:350:G:H2'	23:BB:351:C:O4'	2.22	0.40
23:BB:458:G:O2'	23:BB:459:U:P	2.80	0.40
23:BB:648:G:O2'	23:BB:649:G:H5'	2.22	0.40
23:BB:871:U:H4'	38:BM:68:PHE:CE2	2.57	0.40
23:BB:927:A:H2'	23:BB:928:A:O4'	2.22	0.40
23:BB:948:C:H1'	23:BB:984:A:N3	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:237:ARG:O	25:BC:238:ASN:HB2	2.22	0.40
26:BD:146:ILE:HG13	26:BD:147:GLY:H	1.86	0.40
29:BE:153:LEU:HD12	29:BE:154:ASP:N	2.36	0.40
29:BE:31:VAL:HG21	29:BE:104:ALA:HB2	2.04	0.40
29:BE:98:LYS:CE	29:BE:99:LYS:HG2	2.51	0.40
47:BF:11:VAL:O	47:BF:13:LYS:N	2.54	0.40
48:BG:97:VAL:CB	48:BG:124:CYS:HB2	2.52	0.40
40:BH:62:LEU:HD13	40:BH:66:ASN:HB3	2.04	0.40
41:BJ:101:ILE:O	41:BJ:104:ALA:HB3	2.21	0.40
41:BJ:88:THR:HG23	41:BJ:91:GLU:HB2	2.03	0.40
27:BK:43:ILE:HG22	27:BK:46:ALA:HB2	2.04	0.40
37:BL:121:THR:CG2	37:BL:141:LYS:HB3	2.52	0.40
37:BL:90:VAL:HB	37:BL:122:VAL:CG1	2.47	0.40
37:BL:91:ASP:HB2	37:BL:94:THR:HG23	2.03	0.40
42:BN:82:GLU:C	42:BN:84:GLY:H	2.25	0.40
43:BO:100:HIS:HA	43:BO:104:GLN:HG3	2.03	0.40
28:BP:104:GLY:O	28:BP:106:ALA:N	2.55	0.40
44:BQ:87:VAL:HG12	44:BQ:88:GLU:N	2.36	0.40
45:BS:89:ALA:O	45:BS:90:LYS:HB2	2.22	0.40
50:BT:2:ILE:N	50:BT:2:ILE:HD13	2.36	0.40
39:BX:10:SER:O	39:BX:11:VAL:C	2.59	0.40
39:BX:27:ASN:HD22	39:BX:27:ASN:HA	1.68	0.40
30:BY:9:THR:HB	30:BY:10:ARG:H	1.78	0.40
51:BZ:53:ALA:O	51:BZ:55:GLY:N	2.46	0.40
1:CA:1015:G:H2'	1:CA:1016:A:C8	2.56	0.40
1:CA:1261:A:N7	1:CA:1274:A:H2	2.20	0.40
1:CA:1258:G:N3	1:CA:1278:G:N2	2.70	0.40
1:CA:1456:A:H2'	1:CA:1457:G:O4'	2.21	0.40
1:CA:389:A:H2'	1:CA:389:A:N3	2.36	0.40
1:CA:528:C:H41	11:CL:45:ASN:CG	2.25	0.40
1:CA:529:G:O6	11:CL:45:ASN:HA	2.22	0.40
1:CA:54:C:H2'	1:CA:352:C:H41	1.86	0.40
1:CA:573:A:H2'	1:CA:574:A:C8	2.57	0.40
1:CA:659:U:O2'	1:CA:660:C:H5'	2.22	0.40
18:CB:97:GLY:HA2	18:CB:174:GLU:OE2	2.22	0.40
18:CB:214:GLY:C	18:CB:216:VAL:N	2.73	0.40
3:CD:24:VAL:CG1	3:CD:160:LEU:HB3	2.51	0.40
5:CF:89:VAL:HG13	5:CF:89:VAL:O	2.21	0.40
10:CK:126:ARG:HE	10:CK:126:ARG:CA	2.34	0.40
12:CM:29:SER:CA	12:CM:32:ILE:HG22	2.43	0.40
21:CN:63:CYS:SG	21:CN:64:ARG:N	2.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CR:34:GLU:N	15:CR:34:GLU:OE1	2.50	0.40
16:CS:77:ARG:HB2	16:CS:78:THR:H	1.77	0.40
17:CT:61:ALA:CA	17:CT:67:HIS:H	2.34	0.40
32:D4:25:VAL:O	32:D4:26:ILE:HD13	2.22	0.40
32:D4:27:CYS:HB3	32:D4:33:HIS:HB2	2.03	0.40
32:D4:9:LYS:HD3	32:D4:9:LYS:C	2.42	0.40
23:DB:1104:C:O5'	23:DB:1104:C:H6	2.04	0.40
23:DB:1147:A:O2'	23:DB:1148:U:H5'	2.21	0.40
23:DB:1292:G:H2'	23:DB:1293:C:H6	1.86	0.40
23:DB:138:U:O2'	23:DB:140:C:C5'	2.70	0.40
23:DB:1662:U:O2	23:DB:2687:U:H4'	2.21	0.40
23:DB:179:C:H2'	23:DB:180:G:O4'	2.22	0.40
23:DB:216:A:H2'	23:DB:217:A:C8	2.55	0.40
23:DB:2222:C:H4'	25:DC:184:GLU:OE2	2.22	0.40
23:DB:2391:G:HO2'	23:DB:2392:A:P	2.43	0.40
23:DB:2606:C:O2'	23:DB:2607:G:H5'	2.21	0.40
23:DB:2758:A:O2'	23:DB:2759:G:H5'	2.22	0.40
23:DB:2772:C:H2'	23:DB:2773:C:H6	1.87	0.40
23:DB:616:A:H4'	29:DE:101:TYR:OH	2.21	0.40
23:DB:714:U:H3	23:DB:716:A:H3'	1.86	0.40
23:DB:695:G:C6	23:DB:768:G:C6	3.10	0.40
23:DB:876:C:H5'	23:DB:877:A:O5'	2.21	0.40
47:DF:119:LYS:HA	47:DF:121:PHE:CE2	2.55	0.40
47:DF:121:PHE:CD2	47:DF:121:PHE:N	2.89	0.40
23:DB:1060:U:O4	24:DI:131:THR:HG22	2.20	0.40
27:DK:25:LEU:HD12	27:DK:38:ILE:HB	2.03	0.40
28:DP:111:GLU:CD	28:DP:111:GLU:N	2.75	0.40
44:DQ:33:VAL:O	44:DQ:34:ALA:C	2.60	0.40
44:DQ:108:LEU:HG	49:DR:48:LYS:HG2	2.02	0.40
49:DR:60:LYS:O	49:DR:98:ILE:HA	2.21	0.40
50:DT:32:LEU:H	50:DT:83:ALA:CB	2.34	0.40
50:DT:74:ILE:HG13	50:DT:75:GLY:H	1.87	0.40
50:DT:87:LEU:O	50:DT:88:LYS:O	2.39	0.40
46:DU:18:LYS:O	46:DU:20:LYS:N	2.54	0.40
35:DV:75:GLN:CB	35:DV:90:ASP:HB2	2.52	0.40
52:DW:37:VAL:C	52:DW:38:ARG:HG2	2.42	0.40
52:DW:61:LYS:HB3	52:DW:62:ALA:H	1.49	0.40
39:DX:22:LEU:O	39:DX:22:LEU:HD23	2.22	0.40

There are no symmetry-related clashes.

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	AC	204/232 (88%)	112 (55%)	56 (28%)	36 (18%)	0	2
2	CC	204/232 (88%)	134 (66%)	48 (24%)	22 (11%)	0	8
3	AD	203/205 (99%)	133 (66%)	58 (29%)	12 (6%)	1	19
3	CD	203/205 (99%)	132 (65%)	58 (29%)	13 (6%)	1	18
4	AE	148/166 (89%)	109 (74%)	30 (20%)	9 (6%)	1	18
4	CE	148/166 (89%)	108 (73%)	31 (21%)	9 (6%)	1	18
5	AF	98/135 (73%)	62 (63%)	27 (28%)	9 (9%)	1	12
5	CF	98/135 (73%)	64 (65%)	25 (26%)	9 (9%)	1	12
6	AG	148/178 (83%)	98 (66%)	44 (30%)	6 (4%)	3	25
6	CG	150/178 (84%)	101 (67%)	36 (24%)	13 (9%)	1	12
7	AH	127/129 (98%)	86 (68%)	35 (28%)	6 (5%)	2	23
7	CH	127/129 (98%)	85 (67%)	36 (28%)	6 (5%)	2	23
8	AI	125/129 (97%)	84 (67%)	25 (20%)	16 (13%)	0	5
8	CI	125/129 (97%)	89 (71%)	30 (24%)	6 (5%)	2	23
9	AJ	96/103 (93%)	61 (64%)	18 (19%)	17 (18%)	0	2
9	CJ	96/103 (93%)	62 (65%)	21 (22%)	13 (14%)	0	4
10	AK	115/128 (90%)	85 (74%)	26 (23%)	4 (4%)	3	29
10	CK	115/128 (90%)	84 (73%)	25 (22%)	6 (5%)	2	21
11	AL	121/123 (98%)	71 (59%)	34 (28%)	16 (13%)	0	4
11	CL	121/123 (98%)	72 (60%)	33 (27%)	16 (13%)	0	4
12	AM	112/117 (96%)	69 (62%)	36 (32%)	7 (6%)	1	18
12	CM	111/117 (95%)	77 (69%)	23 (21%)	11 (10%)	0	9
13	AP	80/82 (98%)	53 (66%)	18 (22%)	9 (11%)	0	7
13	CP	78/82 (95%)	53 (68%)	19 (24%)	6 (8%)	1	15
14	AQ	78/83 (94%)	61 (78%)	14 (18%)	3 (4%)	3	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
14	CQ	79/83 (95%)	62 (78%)	15 (19%)	2 (2%)	5	35
15	AR	53/74 (72%)	33 (62%)	17 (32%)	3 (6%)	1	19
15	CR	53/74 (72%)	33 (62%)	16 (30%)	4 (8%)	1	15
16	AS	77/91 (85%)	49 (64%)	21 (27%)	7 (9%)	1	12
16	CS	78/91 (86%)	51 (65%)	20 (26%)	7 (9%)	1	12
17	AT	83/86 (96%)	62 (75%)	16 (19%)	5 (6%)	1	19
17	CT	83/86 (96%)	63 (76%)	14 (17%)	6 (7%)	1	16
18	AB	216/240 (90%)	140 (65%)	53 (24%)	23 (11%)	0	8
18	CB	216/240 (90%)	135 (62%)	59 (27%)	22 (10%)	0	9
19	AU	49/70 (70%)	29 (59%)	13 (26%)	7 (14%)	0	3
19	CU	49/70 (70%)	29 (59%)	15 (31%)	5 (10%)	0	9
20	AO	86/89 (97%)	55 (64%)	24 (28%)	7 (8%)	1	14
20	CO	86/89 (97%)	50 (58%)	29 (34%)	7 (8%)	1	14
21	AN	92/100 (92%)	54 (59%)	29 (32%)	9 (10%)	0	10
21	CN	92/100 (92%)	45 (49%)	31 (34%)	16 (17%)	0	2
24	BI	139/141 (99%)	118 (85%)	16 (12%)	5 (4%)	3	28
24	DI	139/141 (99%)	115 (83%)	19 (14%)	5 (4%)	3	28
25	BC	269/272 (99%)	149 (55%)	68 (25%)	52 (19%)	0	2
25	DC	269/272 (99%)	147 (55%)	70 (26%)	52 (19%)	0	2
26	BD	207/209 (99%)	113 (55%)	58 (28%)	36 (17%)	0	2
26	DD	207/209 (99%)	114 (55%)	58 (28%)	35 (17%)	0	3
27	BK	119/123 (97%)	73 (61%)	24 (20%)	22 (18%)	0	2
27	DK	119/123 (97%)	73 (61%)	25 (21%)	21 (18%)	0	2
28	BP	112/114 (98%)	67 (60%)	28 (25%)	17 (15%)	0	3
28	DP	112/114 (98%)	66 (59%)	31 (28%)	15 (13%)	0	4
29	BE	199/201 (99%)	120 (60%)	49 (25%)	30 (15%)	0	3
29	DE	199/201 (99%)	123 (62%)	47 (24%)	29 (15%)	0	3
30	BY	56/58 (97%)	36 (64%)	16 (29%)	4 (7%)	1	16
30	DY	56/58 (97%)	36 (64%)	14 (25%)	6 (11%)	0	8
31	B0	54/56 (96%)	34 (63%)	10 (18%)	10 (18%)	0	2
31	D0	54/56 (96%)	35 (65%)	9 (17%)	10 (18%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
32	B4	36/38 (95%)	19 (53%)	13 (36%)	4 (11%)	0	7
32	D4	36/38 (95%)	19 (53%)	13 (36%)	4 (11%)	0	7
33	B1	48/54 (89%)	36 (75%)	8 (17%)	4 (8%)	1	13
33	D1	48/54 (89%)	35 (73%)	9 (19%)	4 (8%)	1	13
34	B3	62/64 (97%)	34 (55%)	20 (32%)	8 (13%)	0	5
34	D3	62/64 (97%)	35 (56%)	19 (31%)	8 (13%)	0	5
35	BV	92/94 (98%)	60 (65%)	25 (27%)	7 (8%)	1	15
35	DV	92/94 (98%)	61 (66%)	24 (26%)	7 (8%)	1	15
36	B2	44/46 (96%)	23 (52%)	16 (36%)	5 (11%)	0	6
36	D2	44/46 (96%)	23 (52%)	12 (27%)	9 (20%)	0	2
37	BL	141/144 (98%)	76 (54%)	37 (26%)	28 (20%)	0	2
37	DL	141/144 (98%)	76 (54%)	39 (28%)	26 (18%)	0	2
38	BM	134/136 (98%)	79 (59%)	39 (29%)	16 (12%)	0	6
38	DM	134/136 (98%)	82 (61%)	35 (26%)	17 (13%)	0	5
39	BX	61/63 (97%)	35 (57%)	20 (33%)	6 (10%)	0	10
39	DX	61/63 (97%)	35 (57%)	20 (33%)	6 (10%)	0	10
40	BH	147/149 (99%)	78 (53%)	42 (29%)	27 (18%)	0	2
40	DH	147/149 (99%)	91 (62%)	30 (20%)	26 (18%)	0	2
41	BJ	140/142 (99%)	85 (61%)	37 (26%)	18 (13%)	0	5
41	DJ	140/142 (99%)	85 (61%)	36 (26%)	19 (14%)	0	4
42	BN	118/127 (93%)	73 (62%)	33 (28%)	12 (10%)	0	9
42	DN	118/127 (93%)	74 (63%)	33 (28%)	11 (9%)	0	11
43	BO	114/117 (97%)	68 (60%)	28 (25%)	18 (16%)	0	3
43	DO	114/117 (97%)	66 (58%)	30 (26%)	18 (16%)	0	3
44	BQ	115/117 (98%)	76 (66%)	29 (25%)	10 (9%)	1	12
44	DQ	115/117 (98%)	76 (66%)	30 (26%)	9 (8%)	1	15
45	BS	108/110 (98%)	59 (55%)	34 (32%)	15 (14%)	0	4
45	DS	108/110 (98%)	60 (56%)	33 (31%)	15 (14%)	0	4
46	BU	100/103 (97%)	58 (58%)	25 (25%)	17 (17%)	0	3
46	DU	100/103 (97%)	57 (57%)	24 (24%)	19 (19%)	0	2
47	BF	176/178 (99%)	91 (52%)	51 (29%)	34 (19%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
47	DF	176/178 (99%)	93 (53%)	49 (28%)	34 (19%)	0	2
48	BG	174/176 (99%)	100 (58%)	42 (24%)	32 (18%)	0	2
48	DG	174/176 (99%)	101 (58%)	42 (24%)	31 (18%)	0	2
49	BR	101/103 (98%)	57 (56%)	26 (26%)	18 (18%)	0	2
49	DR	101/103 (98%)	58 (57%)	26 (26%)	17 (17%)	0	3
50	BT	91/100 (91%)	40 (44%)	40 (44%)	11 (12%)	0	6
50	DT	91/100 (91%)	41 (45%)	39 (43%)	11 (12%)	0	6
51	BZ	75/78 (96%)	53 (71%)	16 (21%)	6 (8%)	1	14
51	DZ	75/78 (96%)	54 (72%)	14 (19%)	7 (9%)	0	11
52	BW	77/84 (92%)	29 (38%)	24 (31%)	24 (31%)	0	0
52	DW	77/84 (92%)	27 (35%)	26 (34%)	24 (31%)	0	0
All	All	11241/11914 (94%)	6932 (62%)	2908 (26%)	1401 (12%)	0	5

All (1401) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	AC	14	VAL
2	AC	19	SER
2	AC	26	LYS
2	AC	47	ALA
2	AC	54	ILE
2	AC	81	GLU
2	AC	100	ILE
2	AC	127	VAL
2	AC	130	ARG
2	AC	205	GLU
4	AE	20	VAL
6	AG	15	PRO
6	AG	112	ASP
7	AH	66	GLN
8	AI	8	THR
8	AI	71	ILE
8	AI	127	SER
9	AJ	57	VAL
9	AJ	99	GLN
11	AL	42	LYS
12	AM	14	ALA
12	AM	49	GLU

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Mol	Chain	Res	Type
12	AM	105	ALA
13	AP	44	SER
16	AS	66	VAL
18	AB	14	HIS
18	AB	15	PHE
18	AB	19	THR
18	AB	22	TRP
18	AB	41	ASN
18	AB	76	SER
18	AB	81	ASP
18	AB	209	VAL
19	AU	35	GLU
20	AO	21	ASP
21	AN	50	LEU
24	BI	18	ASN
25	BC	4	LYS
25	BC	51	ARG
25	BC	65	ASP
25	BC	68	ARG
25	BC	117	SER
25	BC	123	ILE
25	BC	135	PRO
25	BC	141	HIS
25	BC	145	MET
25	BC	202	ARG
25	BC	203	VAL
25	BC	254	LYS
26	BD	9	VAL
26	BD	74	GLU
26	BD	91	THR
26	BD	107	VAL
26	BD	122	VAL
26	BD	143	PRO
26	BD	170	VAL
26	BD	184	ARG
26	BD	189	VAL
27	BK	18	ARG
27	BK	35	VAL
27	BK	43	ILE
27	BK	72	PRO
27	BK	89	ASN
27	BK	120	PRO

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Mol	Chain	Res	Type
28	BP	25	VAL
28	BP	37	LYS
28	BP	50	ARG
28	BP	75	THR
28	BP	100	ARG
29	BE	7	ASP
29	BE	14	VAL
29	BE	45	ALA
29	BE	70	SER
29	BE	79	ARG
30	BY	2	LYS
31	B0	23	ALA
31	B0	42	ILE
31	B0	51	ARG
32	B4	37	GLN
33	B1	4	ILE
34	B3	29	ARG
34	B3	49	VAL
35	BV	75	GLN
37	BL	9	ALA
37	BL	81	ASP
37	BL	89	VAL
37	BL	94	THR
37	BL	100	ILE
37	BL	116	VAL
37	BL	143	GLU
38	BM	24	THR
38	BM	59	ARG
40	BH	3	VAL
40	BH	14	SER
40	BH	28	ASN
40	BH	32	PRO
40	BH	33	GLN
40	BH	122	LEU
41	BJ	4	PHE
41	BJ	41	LYS
41	BJ	44	TYR
41	BJ	45	THR
41	BJ	64	VAL
41	BJ	73	VAL
41	BJ	81	ILE
42	BN	11	ASN

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Mol	Chain	Res	Type
42	BN	13	ASN
43	BO	6	ALA
43	BO	8	ILE
43	BO	113	ALA
44	BQ	30	VAL
44	BQ	86	SER
45	BS	13	SER
45	BS	27	LYS
45	BS	61	ASN
46	BU	18	LYS
46	BU	42	LYS
46	BU	49	PRO
47	BF	11	VAL
47	BF	32	LYS
47	BF	70	ARG
47	BF	112	ASP
47	BF	135	ILE
47	BF	138	PRO
47	BF	149	ARG
48	BG	11	PRO
48	BG	83	THR
48	BG	85	LYS
48	BG	94	ARG
48	BG	117	PRO
48	BG	157	LYS
48	BG	170	THR
49	BR	53	PHE
49	BR	98	ILE
49	BR	99	THR
50	BT	16	VAL
50	BT	58	VAL
50	BT	88	LYS
51	BZ	33	LEU
51	BZ	70	GLU
51	BZ	77	LYS
52	BW	9	THR
52	BW	30	VAL
52	BW	50	VAL
52	BW	60	ALA
52	BW	61	LYS
52	BW	62	ALA
52	BW	77	LYS

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Mol	Chain	Res	Type
4	CE	20	VAL
7	CH	66	GLN
11	CL	42	LYS
13	CP	44	SER
19	CU	35	GLU
2	CC	19	SER
2	CC	54	ILE
2	CC	83	VAL
2	CC	100	ILE
2	CC	205	GLU
6	CG	3	ARG
6	CG	71	THR
8	CI	8	THR
8	CI	57	VAL
9	CJ	37	ARG
9	CJ	57	VAL
12	CM	46	GLU
12	CM	105	ALA
12	CM	111	PRO
21	CN	29	ILE
21	CN	50	LEU
21	CN	61	ASN
21	CN	75	LYS
21	CN	80	ARG
18	CB	15	PHE
18	CB	22	TRP
18	CB	49	PHE
18	CB	94	ARG
18	CB	121	GLN
18	CB	141	GLU
20	CO	29	VAL
24	DI	5	GLN
24	DI	18	ASN
25	DC	4	LYS
25	DC	51	ARG
25	DC	65	ASP
25	DC	68	ARG
25	DC	117	SER
25	DC	123	ILE
25	DC	135	PRO
25	DC	141	HIS
25	DC	145	MET

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Mol	Chain	Res	Type
25	DC	202	ARG
25	DC	203	VAL
25	DC	254	LYS
26	DD	9	VAL
26	DD	74	GLU
26	DD	91	THR
26	DD	107	VAL
26	DD	122	VAL
26	DD	143	PRO
26	DD	170	VAL
26	DD	184	ARG
26	DD	189	VAL
27	DK	18	ARG
27	DK	35	VAL
27	DK	43	ILE
27	DK	72	PRO
27	DK	89	ASN
27	DK	120	PRO
28	DP	25	VAL
28	DP	37	LYS
28	DP	50	ARG
28	DP	75	THR
28	DP	100	ARG
29	DE	7	ASP
29	DE	14	VAL
29	DE	45	ALA
29	DE	70	SER
30	DY	2	LYS
31	D0	17	SER
31	D0	23	ALA
31	D0	42	ILE
31	D0	51	ARG
32	D4	37	GLN
33	D1	4	ILE
34	D3	29	ARG
34	D3	49	VAL
35	DV	75	GLN
37	DL	9	ALA
37	DL	81	ASP
37	DL	89	VAL
37	DL	94	THR
37	DL	100	ILE

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Mol	Chain	Res	Type
37	DL	116	VAL
37	DL	143	GLU
38	DM	24	THR
38	DM	59	ARG
40	DH	3	VAL
40	DH	14	SER
40	DH	28	ASN
40	DH	32	PRO
40	DH	33	GLN
40	DH	110	VAL
40	DH	114	GLU
40	DH	148	ALA
41	DJ	4	PHE
41	DJ	41	LYS
41	DJ	44	TYR
41	DJ	45	THR
41	DJ	64	VAL
41	DJ	73	VAL
41	DJ	81	ILE
42	DN	11	ASN
42	DN	13	ASN
43	DO	6	ALA
43	DO	8	ILE
43	DO	113	ALA
44	DQ	30	VAL
44	DQ	86	SER
45	DS	13	SER
45	DS	27	LYS
45	DS	42	LYS
45	DS	61	ASN
46	DU	18	LYS
46	DU	42	LYS
46	DU	49	PRO
47	DF	11	VAL
47	DF	32	LYS
47	DF	70	ARG
47	DF	112	ASP
47	DF	135	ILE
47	DF	138	PRO
47	DF	148	VAL
47	DF	149	ARG
48	DG	9	VAL

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Mol	Chain	Res	Type
48	DG	11	PRO
48	DG	85	LYS
48	DG	94	ARG
48	DG	117	PRO
48	DG	157	LYS
49	DR	53	PHE
49	DR	98	ILE
49	DR	99	THR
50	DT	16	VAL
50	DT	58	VAL
50	DT	88	LYS
51	DZ	33	LEU
51	DZ	70	GLU
51	DZ	77	LYS
52	DW	9	THR
52	DW	30	VAL
52	DW	32	ALA
52	DW	50	VAL
52	DW	60	ALA
52	DW	61	LYS
52	DW	62	ALA
52	DW	77	LYS
2	AC	25	THR
2	AC	59	PRO
2	AC	83	VAL
2	AC	105	VAL
2	AC	116	ALA
2	AC	167	TYR
2	AC	180	ASP
3	AD	22	SER
3	AD	175	GLY
3	AD	178	GLU
4	AE	71	ILE
5	AF	48	ALA
5	AF	54	LEU
5	AF	62	MET
5	AF	82	ASP
5	AF	92	THR
8	AI	55	ASP
8	AI	106	ASP
8	AI	120	ALA
9	AJ	34	ALA

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Mol	Chain	Res	Type
9	AJ	36	VAL
9	AJ	51	VAL
9	AJ	56	HIS
10	AK	88	PRO
11	AL	13	ARG
11	AL	14	LYS
11	AL	23	LEU
11	AL	24	GLU
11	AL	71	HIS
11	AL	117	GLY
11	AL	122	LYS
12	AM	70	ARG
13	AP	49	GLY
13	AP	79	ASN
15	AR	43	ILE
16	AS	5	LYS
16	AS	77	ARG
17	AT	42	ASP
17	AT	70	LYS
17	AT	85	LEU
18	AB	18	GLN
18	AB	64	GLY
18	AB	94	ARG
18	AB	188	THR
18	AB	204	ASP
19	AU	34	ARG
19	AU	36	PHE
20	AO	34	ALA
20	AO	37	ASN
20	AO	76	ALA
21	AN	62	ARG
24	BI	14	ALA
24	BI	23	VAL
24	BI	64	ARG
25	BC	3	VAL
25	BC	34	GLU
25	BC	36	ASN
25	BC	37	SER
25	BC	52	HIS
25	BC	59	GLN
25	BC	62	ARG
25	BC	63	ILE

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Mol	Chain	Res	Type
25	BC	64	VAL
25	BC	70	LYS
25	BC	121	ALA
25	BC	140	VAL
25	BC	151	GLY
25	BC	189	ALA
25	BC	190	THR
25	BC	222	THR
25	BC	239	PHE
26	BD	10	GLY
26	BD	11	MET
26	BD	53	GLY
26	BD	93	GLY
26	BD	102	ALA
26	BD	118	PHE
26	BD	119	ALA
26	BD	121	THR
26	BD	145	SER
26	BD	159	LYS
26	BD	169	ARG
26	BD	174	SER
26	BD	182	ALA
27	BK	31	ARG
27	BK	110	GLU
28	BP	101	GLU
29	BE	12	LEU
29	BE	27	LEU
29	BE	42	GLY
29	BE	63	LYS
29	BE	81	GLY
29	BE	86	ALA
29	BE	166	LYS
30	BY	39	ASP
31	B0	17	SER
31	B0	54	ILE
32	B4	4	ARG
33	B1	51	ALA
34	B3	31	ILE
34	B3	58	ILE
35	BV	71	LYS
36	B2	5	PHE
36	B2	45	SER

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Mol	Chain	Res	Type
37	BL	15	ALA
37	BL	24	GLY
37	BL	28	GLY
37	BL	51	GLU
37	BL	62	PRO
37	BL	99	ASN
37	BL	111	ILE
38	BM	20	LEU
38	BM	69	PRO
39	BX	58	ASN
40	BH	7	ASP
40	BH	27	ARG
40	BH	67	ALA
40	BH	73	ASN
40	BH	89	LYS
40	BH	97	ARG
40	BH	125	THR
40	BH	144	VAL
41	BJ	5	THR
41	BJ	43	GLU
41	BJ	84	ILE
41	BJ	124	VAL
42	BN	68	ALA
42	BN	93	GLY
42	BN	100	CYS
43	BO	99	TYR
44	BQ	87	VAL
44	BQ	89	ILE
45	BS	3	THR
45	BS	25	ARG
45	BS	28	LYS
45	BS	41	LYS
45	BS	42	LYS
45	BS	96	ILE
46	BU	12	VAL
46	BU	19	GLY
46	BU	50	ALA
46	BU	92	VAL
47	BF	80	GLN
47	BF	87	LYS
47	BF	92	GLY
47	BF	136	ILE

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Mol	Chain	Res	Type
47	BF	148	VAL
48	BG	9	VAL
48	BG	84	LYS
48	BG	163	TYR
49	BR	7	SER
49	BR	16	GLU
49	BR	55	ASP
49	BR	79	ARG
50	BT	38	ALA
50	BT	39	THR
51	BZ	71	LEU
52	BW	14	ASP
52	BW	32	ALA
52	BW	34	SER
52	BW	36	ILE
52	BW	59	PHE
3	CD	22	SER
3	CD	175	GLY
3	CD	178	GLU
3	CD	192	ALA
4	CE	71	ILE
5	CF	48	ALA
5	CF	54	LEU
5	CF	62	MET
5	CF	92	THR
5	CF	95	ALA
10	CK	88	PRO
11	CL	13	ARG
11	CL	14	LYS
11	CL	23	LEU
11	CL	24	GLU
11	CL	71	HIS
11	CL	117	GLY
11	CL	122	LYS
13	CP	28	ARG
13	CP	49	GLY
15	CR	43	ILE
17	CT	42	ASP
17	CT	70	LYS
17	CT	85	LEU
19	CU	34	ARG
19	CU	36	PHE

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Mol	Chain	Res	Type
2	CC	25	THR
2	CC	26	LYS
2	CC	59	PRO
2	CC	72	PRO
2	CC	80	GLY
2	CC	94	ALA
2	CC	112	ALA
2	CC	166	TRP
2	CC	180	ASP
2	CC	192	TYR
6	CG	30	MET
6	CG	130	LYS
8	CI	91	GLU
9	CJ	34	ALA
9	CJ	36	VAL
9	CJ	61	ALA
9	CJ	92	LEU
12	CM	3	ILE
12	CM	6	ILE
21	CN	20	PHE
21	CN	27	LYS
21	CN	57	SER
16	CS	65	MET
16	CS	77	ARG
18	CB	9	LEU
18	CB	65	LYS
18	CB	87	ASP
18	CB	100	LEU
18	CB	163	ILE
20	CO	8	THR
20	CO	48	LYS
25	DC	3	VAL
25	DC	34	GLU
25	DC	36	ASN
25	DC	37	SER
25	DC	52	HIS
25	DC	59	GLN
25	DC	62	ARG
25	DC	63	ILE
25	DC	64	VAL
25	DC	121	ALA
25	DC	140	VAL

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Mol	Chain	Res	Type
25	DC	151	GLY
25	DC	189	ALA
25	DC	190	THR
25	DC	222	THR
25	DC	239	PHE
26	DD	10	GLY
26	DD	11	MET
26	DD	53	GLY
26	DD	93	GLY
26	DD	102	ALA
26	DD	118	PHE
26	DD	119	ALA
26	DD	121	THR
26	DD	145	SER
26	DD	159	LYS
26	DD	169	ARG
26	DD	174	SER
26	DD	182	ALA
26	DD	197	THR
27	DK	31	ARG
27	DK	110	GLU
28	DP	86	LYS
28	DP	101	GLU
29	DE	12	LEU
29	DE	27	LEU
29	DE	42	GLY
29	DE	63	LYS
29	DE	79	ARG
29	DE	86	ALA
29	DE	166	LYS
30	DY	39	ASP
32	D4	4	ARG
33	D1	51	ALA
34	D3	31	ILE
34	D3	58	ILE
35	DV	71	LYS
36	D2	5	PHE
36	D2	45	SER
37	DL	15	ALA
37	DL	24	GLY
37	DL	51	GLU
37	DL	62	PRO

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Mol	Chain	Res	Type
37	DL	99	ASN
37	DL	111	ILE
37	DL	138	ALA
38	DM	20	LEU
38	DM	69	PRO
39	DX	58	ASN
40	DH	7	ASP
40	DH	27	ARG
40	DH	40	THR
40	DH	85	GLY
40	DH	96	THR
41	DJ	3	THR
41	DJ	5	THR
41	DJ	43	GLU
41	DJ	84	ILE
42	DN	68	ALA
42	DN	93	GLY
42	DN	100	CYS
42	DN	101	GLY
43	DO	45	SER
43	DO	68	LYS
43	DO	99	TYR
43	DO	115	LEU
44	DQ	87	VAL
44	DQ	89	ILE
45	DS	3	THR
45	DS	25	ARG
45	DS	28	LYS
45	DS	41	LYS
45	DS	96	ILE
46	DU	12	VAL
46	DU	19	GLY
46	DU	50	ALA
46	DU	92	VAL
47	DF	42	ALA
47	DF	80	GLN
47	DF	87	LYS
47	DF	92	GLY
47	DF	136	ILE
48	DG	83	THR
48	DG	84	LYS
48	DG	163	TYR

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Mol	Chain	Res	Type
48	DG	170	THR
49	DR	7	SER
49	DR	16	GLU
49	DR	55	ASP
49	DR	70	GLU
49	DR	79	ARG
50	DT	38	ALA
50	DT	39	THR
51	DZ	35	SER
51	DZ	71	LEU
52	DW	14	ASP
52	DW	23	LYS
52	DW	29	SER
52	DW	34	SER
52	DW	36	ILE
52	DW	59	PHE
2	AC	104	GLU
2	AC	107	LYS
2	AC	113	LYS
2	AC	166	TRP
2	AC	178	ARG
2	AC	190	THR
3	AD	26	ALA
3	AD	27	ILE
3	AD	192	ALA
4	AE	146	MET
5	AF	85	ILE
5	AF	95	ALA
5	AF	98	GLU
6	AG	23	ALA
7	AH	72	GLU
8	AI	50	PRO
8	AI	82	ILE
8	AI	99	LYS
9	AJ	62	ARG
9	AJ	93	ALA
10	AK	14	GLN
11	AL	56	LEU
11	AL	72	ASN
12	AM	68	LEU
12	AM	98	GLY
12	AM	104	ASN

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Mol	Chain	Res	Type
13	AP	28	ARG
13	AP	52	LEU
14	AQ	16	MET
18	AB	75	ALA
18	AB	86	CYS
18	AB	99	MET
18	AB	128	LEU
19	AU	9	GLU
19	AU	37	TYR
20	AO	20	ASN
20	AO	36	ILE
21	AN	2	LYS
21	AN	27	LYS
21	AN	70	HIS
25	BC	15	VAL
25	BC	88	ALA
25	BC	109	LEU
25	BC	200	MET
25	BC	233	GLY
26	BD	31	ALA
26	BD	77	ARG
26	BD	106	LYS
26	BD	136	ASN
26	BD	167	ASN
26	BD	181	ASP
26	BD	195	GLY
26	BD	197	THR
27	BK	14	SER
27	BK	17	ARG
27	BK	30	ARG
27	BK	73	ASP
27	BK	90	ASN
28	BP	36	LYS
28	BP	86	LYS
29	BE	52	VAL
29	BE	123	LYS
29	BE	153	LEU
29	BE	188	MET
31	B0	45	ASP
32	B4	20	ASP
33	B1	50	GLU
35	BV	6	ALA

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Mol	Chain	Res	Type
37	BL	5	THR
37	BL	29	LYS
37	BL	36	LYS
37	BL	58	TYR
37	BL	138	ALA
38	BM	13	HIS
38	BM	43	ALA
38	BM	83	GLY
38	BM	134	THR
39	BX	9	LYS
39	BX	37	LEU
39	BX	45	GLN
40	BH	6	LEU
40	BH	11	ASN
40	BH	29	PHE
40	BH	31	VAL
40	BH	54	LEU
40	BH	83	LYS
40	BH	99	ILE
40	BH	131	SER
40	BH	140	ALA
41	BJ	2	LYS
41	BJ	3	THR
42	BN	71	ARG
42	BN	72	ASP
42	BN	98	LEU
42	BN	101	GLY
43	BO	45	SER
43	BO	51	ALA
43	BO	53	THR
43	BO	66	GLY
43	BO	68	LYS
43	BO	100	HIS
43	BO	115	LEU
44	BQ	91	ARG
45	BS	21	ALA
45	BS	46	LEU
46	BU	6	ARG
46	BU	82	VAL
46	BU	85	ARG
46	BU	101	THR
47	BF	18	GLU

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Mol	Chain	Res	Type
47	BF	42	ALA
47	BF	78	ILE
47	BF	93	GLU
47	BF	106	ALA
47	BF	122	ASP
47	BF	141	ASP
48	BG	2	ARG
48	BG	32	LEU
48	BG	61	TRP
48	BG	152	ARG
48	BG	164	ALA
49	BR	24	LYS
49	BR	43	ASN
49	BR	70	GLU
49	BR	83	TYR
51	BZ	35	SER
52	BW	23	LYS
52	BW	29	SER
52	BW	53	GLY
52	BW	69	GLU
52	BW	70	VAL
3	CD	26	ALA
3	CD	27	ILE
3	CD	42	ALA
3	CD	191	SER
4	CE	146	MET
5	CF	82	ASP
5	CF	85	ILE
5	CF	98	GLU
7	CH	72	GLU
10	CK	14	GLN
11	CL	56	LEU
11	CL	70	GLY
11	CL	72	ASN
13	CP	52	LEU
17	CT	46	ALA
19	CU	9	GLU
19	CU	37	TYR
2	CC	145	ALA
2	CC	167	TYR
6	CG	70	PRO
6	CG	129	ASN

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Mol	Chain	Res	Type
8	CI	26	LYS
9	CJ	62	ARG
12	CM	69	ARG
12	CM	104	ASN
21	CN	8	ARG
21	CN	62	ARG
16	CS	5	LYS
18	CB	18	GLN
18	CB	59	ILE
18	CB	63	LYS
18	CB	127	LYS
18	CB	128	LEU
18	CB	188	THR
20	CO	25	THR
20	CO	47	LYS
20	CO	50	HIS
24	DI	23	VAL
25	DC	15	VAL
25	DC	70	LYS
25	DC	88	ALA
25	DC	109	LEU
25	DC	200	MET
26	DD	31	ALA
26	DD	77	ARG
26	DD	106	LYS
26	DD	136	ASN
26	DD	167	ASN
26	DD	181	ASP
26	DD	195	GLY
27	DK	14	SER
27	DK	17	ARG
27	DK	30	ARG
27	DK	73	ASP
27	DK	90	ASN
28	DP	36	LYS
29	DE	52	VAL
29	DE	81	GLY
29	DE	84	THR
29	DE	97	ASN
29	DE	123	LYS
29	DE	153	LEU
29	DE	188	MET

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Mol	Chain	Res	Type
31	D0	45	ASP
31	D0	54	ILE
32	D4	16	ILE
32	D4	20	ASP
33	D1	50	GLU
34	D3	18	LYS
35	DV	6	ALA
36	D2	14	ARG
37	DL	5	THR
37	DL	28	GLY
37	DL	29	LYS
37	DL	54	GLN
37	DL	58	TYR
38	DM	43	ALA
38	DM	83	GLY
39	DX	37	LEU
39	DX	45	GLN
40	DH	6	LEU
40	DH	29	PHE
40	DH	31	VAL
40	DH	56	ALA
40	DH	109	GLU
40	DH	111	ALA
40	DH	120	GLY
41	DJ	2	LYS
41	DJ	124	VAL
42	DN	72	ASP
42	DN	98	LEU
43	DO	51	ALA
43	DO	53	THR
43	DO	89	ASP
43	DO	100	HIS
44	DQ	91	ARG
45	DS	21	ALA
45	DS	43	ALA
45	DS	46	LEU
46	DU	6	ARG
46	DU	82	VAL
46	DU	85	ARG
46	DU	101	THR
47	DF	18	GLU
47	DF	78	ILE

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Mol	Chain	Res	Type
47	DF	93	GLU
47	DF	106	ALA
47	DF	110	ILE
47	DF	122	ASP
47	DF	141	ASP
48	DG	2	ARG
48	DG	32	LEU
48	DG	61	TRP
48	DG	152	ARG
48	DG	164	ALA
49	DR	24	LYS
49	DR	43	ASN
49	DR	83	TYR
50	DT	29	THR
52	DW	69	GLU
52	DW	70	VAL
2	AC	3	LYS
2	AC	17	TRP
2	AC	23	ALA
2	AC	50	SER
2	AC	64	ARG
2	AC	109	GLU
2	AC	145	ALA
2	AC	165	GLU
2	AC	179	ALA
3	AD	42	ALA
3	AD	68	GLU
3	AD	191	SER
4	AE	77	ASN
6	AG	10	LYS
6	AG	92	PRO
8	AI	24	ASN
8	AI	25	GLY
8	AI	26	LYS
8	AI	80	HIS
9	AJ	16	ARG
9	AJ	42	LEU
9	AJ	74	VAL
9	AJ	75	ASP
10	AK	71	ASP
11	AL	70	GLY
13	AP	42	ILE

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Mol	Chain	Res	Type
13	AP	81	ALA
16	AS	13	HIS
17	AT	46	ALA
18	AB	73	ARG
18	AB	87	ASP
18	AB	107	ARG
18	AB	126	ASP
18	AB	205	ALA
21	AN	11	LYS
25	BC	29	PHE
25	BC	105	ALA
25	BC	196	ASN
25	BC	204	LEU
25	BC	209	ALA
25	BC	246	PRO
25	BC	250	GLN
26	BD	92	VAL
26	BD	109	VAL
26	BD	206	ALA
27	BK	26	GLY
27	BK	119	ALA
28	BP	31	VAL
28	BP	59	THR
28	BP	65	ASN
28	BP	76	HIS
28	BP	105	LYS
29	BE	6	LYS
29	BE	13	THR
29	BE	84	THR
29	BE	97	ASN
29	BE	183	PHE
30	BY	4	ILE
30	BY	9	THR
31	B0	19	ASP
31	B0	26	SER
32	B4	16	ILE
33	B1	36	LYS
34	B3	18	LYS
34	B3	22	LYS
34	B3	50	SER
35	BV	45	ASP
35	BV	67	GLY

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Mol	Chain	Res	Type
37	BL	3	LEU
37	BL	54	GLN
38	BM	42	THR
38	BM	65	ILE
38	BM	106	ASP
38	BM	116	ALA
39	BX	51	ALA
40	BH	12	LEU
40	BH	77	THR
40	BH	104	THR
41	BJ	25	LEU
41	BJ	47	HIS
41	BJ	128	ASN
43	BO	65	THR
43	BO	89	ASP
44	BQ	10	ARG
44	BQ	27	ARG
45	BS	18	ARG
45	BS	43	ALA
46	BU	16	LYS
46	BU	41	VAL
46	BU	54	PRO
47	BF	28	PRO
47	BF	36	ASN
47	BF	69	ALA
47	BF	110	ILE
47	BF	147	ARG
48	BG	28	LYS
48	BG	29	ASN
48	BG	31	GLU
48	BG	38	ASP
48	BG	93	TYR
48	BG	125	PRO
48	BG	151	ARG
49	BR	42	ALA
49	BR	52	PRO
50	BT	18	GLU
50	BT	29	THR
51	BZ	31	PRO
52	BW	28	GLU
52	BW	40	ARG
52	BW	56	HIS

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Mol	Chain	Res	Type
3	CD	68	GLU
4	CE	77	ASN
10	CK	71	ASP
11	CL	47	ALA
11	CL	74	GLN
13	CP	42	ILE
14	CQ	16	MET
2	CC	146	LYS
8	CI	67	LYS
9	CJ	15	HIS
9	CJ	75	ASP
9	CJ	91	ASP
9	CJ	93	ALA
12	CM	7	ASN
12	CM	20	SER
12	CM	71	GLU
21	CN	63	CYS
21	CN	86	ALA
16	CS	27	LYS
16	CS	53	GLY
18	CB	97	GLY
18	CB	124	THR
20	CO	18	ASP
25	DC	29	PHE
25	DC	105	ALA
25	DC	196	ASN
25	DC	204	LEU
25	DC	233	GLY
25	DC	246	PRO
25	DC	250	GLN
26	DD	92	VAL
26	DD	109	VAL
27	DK	26	GLY
27	DK	113	MET
27	DK	117	SER
27	DK	119	ALA
29	DE	6	LYS
29	DE	83	VAL
29	DE	133	LEU
29	DE	183	PHE
30	DY	4	ILE
30	DY	9	THR

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Mol	Chain	Res	Type
31	D0	19	ASP
31	D0	26	SER
33	D1	36	LYS
34	D3	22	LYS
34	D3	50	SER
35	DV	45	ASP
35	DV	67	GLY
37	DL	3	LEU
37	DL	36	LYS
38	DM	13	HIS
38	DM	42	THR
38	DM	65	ILE
38	DM	106	ASP
38	DM	116	ALA
38	DM	134	THR
39	DX	9	LYS
40	DH	11	ASN
40	DH	86	ASP
41	DJ	25	LEU
41	DJ	47	HIS
42	DN	71	ARG
43	DO	65	THR
43	DO	66	GLY
43	DO	77	ALA
44	DQ	10	ARG
44	DQ	27	ARG
45	DS	18	ARG
46	DU	16	LYS
46	DU	41	VAL
46	DU	51	LEU
46	DU	54	PRO
47	DF	28	PRO
47	DF	36	ASN
47	DF	62	GLN
47	DF	147	ARG
48	DG	28	LYS
48	DG	29	ASN
48	DG	31	GLU
48	DG	38	ASP
48	DG	125	PRO
48	DG	151	ARG
49	DR	42	ALA

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Mol	Chain	Res	Type
50	DT	11	LEU
51	DZ	31	PRO
52	DW	28	GLU
52	DW	40	ARG
52	DW	53	GLY
3	AD	31	CYS
8	AI	108	ARG
11	AL	47	ALA
11	AL	74	GLN
13	AP	46	LYS
16	AS	4	LEU
16	AS	64	GLU
17	AT	66	ILE
19	AU	17	ARG
19	AU	41	THR
21	AN	21	ALA
21	AN	67	GLY
25	BC	11	GLY
27	BK	4	GLU
27	BK	46	ALA
27	BK	113	MET
27	BK	117	SER
28	BP	113	LEU
29	BE	83	VAL
29	BE	133	LEU
29	BE	189	THR
31	B0	48	TYR
34	B3	6	VAL
35	BV	84	PRO
36	B2	14	ARG
36	B2	18	PHE
37	BL	84	LYS
38	BM	60	GLN
38	BM	72	PRO
38	BM	73	ILE
38	BM	95	LEU
41	BJ	129	GLU
41	BJ	134	ALA
42	BN	59	SER
42	BN	104	ALA
43	BO	47	VAL
43	BO	77	ALA

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Mol	Chain	Res	Type
46	BU	51	LEU
47	BF	12	VAL
47	BF	47	LYS
47	BF	132	ARG
47	BF	173	ASP
48	BG	16	VAL
48	BG	96	ALA
48	BG	97	VAL
48	BG	111	PRO
49	BR	101	ILE
50	BT	35	ALA
50	BT	55	VAL
52	BW	74	LYS
3	CD	31	CYS
7	CH	70	VAL
10	CK	125	LYS
17	CT	50	PHE
17	CT	66	ILE
2	CC	3	LYS
2	CC	21	TRP
6	CG	20	GLU
6	CG	64	ALA
6	CG	66	GLU
6	CG	151	ALA
9	CJ	56	HIS
9	CJ	74	VAL
12	CM	47	LEU
21	CN	26	LEU
21	CN	87	ALA
18	CB	99	MET
24	DI	6	ALA
24	DI	14	ALA
25	DC	11	GLY
25	DC	53	ILE
25	DC	209	ALA
25	DC	263	ASP
26	DD	173	GLN
26	DD	206	ALA
27	DK	46	ALA
28	DP	31	VAL
28	DP	59	THR
28	DP	76	HIS

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Mol	Chain	Res	Type
28	DP	105	LYS
29	DE	13	THR
29	DE	69	ARG
29	DE	189	THR
30	DY	51	SER
34	D3	6	VAL
35	DV	84	PRO
36	D2	18	PHE
36	D2	40	ALA
37	DL	41	ARG
37	DL	52	GLY
37	DL	103	ILE
38	DM	72	PRO
38	DM	73	ILE
38	DM	95	LEU
39	DX	51	ALA
40	DH	12	LEU
40	DH	121	VAL
41	DJ	128	ASN
41	DJ	129	GLU
42	DN	59	SER
42	DN	107	ASN
43	DO	47	VAL
44	DQ	39	ILE
46	DU	67	SER
47	DF	12	VAL
47	DF	69	ALA
47	DF	132	ARG
47	DF	145	VAL
47	DF	173	ASP
48	DG	16	VAL
48	DG	93	TYR
48	DG	96	ALA
48	DG	97	VAL
48	DG	111	PRO
48	DG	155	PRO
49	DR	23	GLU
49	DR	52	PRO
49	DR	101	ILE
50	DT	35	ALA
50	DT	55	VAL
50	DT	65	GLY

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Mol	Chain	Res	Type
52	DW	56	HIS
52	DW	74	LYS
2	AC	112	ALA
2	AC	143	LEU
7	AH	70	VAL
7	AH	91	LEU
8	AI	74	GLN
11	AL	84	GLY
13	AP	31	ARG
14	AQ	35	LYS
14	AQ	48	GLU
24	BI	49	GLU
25	BC	53	ILE
25	BC	149	LYS
25	BC	249	VAL
25	BC	251	THR
26	BD	162	ALA
26	BD	173	GLN
27	BK	3	GLN
28	BP	83	ILE
29	BE	69	ARG
29	BE	96	VAL
31	B0	46	GLY
37	BL	19	LEU
37	BL	101	ILE
37	BL	103	ILE
42	BN	10	LEU
44	BQ	32	ARG
44	BQ	39	ILE
45	BS	15	GLN
46	BU	5	ARG
47	BF	81	GLY
47	BF	145	VAL
48	BG	89	VAL
48	BG	112	VAL
48	BG	137	LYS
48	BG	155	PRO
49	BR	65	ALA
50	BT	67	VAL
52	BW	12	GLY
3	CD	182	LYS
4	CE	132	PRO

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Mol	Chain	Res	Type
7	CH	91	LEU
10	CK	124	LYS
13	CP	31	ARG
14	CQ	48	GLU
15	CR	46	THR
2	CC	14	VAL
21	CN	56	PRO
21	CN	66	THR
16	CS	34	SER
25	DC	249	VAL
25	DC	251	THR
27	DK	121	GLU
30	DY	34	THR
31	D0	48	TYR
36	D2	8	SER
36	D2	21	ARG
36	D2	25	LYS
37	DL	101	ILE
38	DM	56	ALA
38	DM	82	MET
40	DH	41	LYS
41	DJ	134	ALA
45	DS	15	GLN
46	DU	5	ARG
47	DF	81	GLY
48	DG	89	VAL
48	DG	112	VAL
50	DT	67	VAL
51	DZ	28	ARG
52	DW	12	GLY
52	DW	33	GLY
3	AD	86	GLY
4	AE	43	GLY
6	AG	54	GLY
11	AL	3	VAL
11	AL	43	LYS
25	BC	232	GLY
29	BE	148	ILE
37	BL	52	GLY
39	BX	11	VAL
40	BH	147	VAL
43	BO	101	GLY

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Mol	Chain	Res	Type
47	BF	84	ILE
48	BG	167	VAL
50	BT	65	GLY
3	CD	86	GLY
4	CE	43	GLY
4	CE	56	PRO
5	CF	89	VAL
7	CH	33	VAL
11	CL	43	LYS
11	CL	84	GLY
8	CI	25	GLY
25	DC	248	GLY
27	DK	93	GLN
28	DP	63	ILE
28	DP	83	ILE
29	DE	96	VAL
36	D2	9	VAL
39	DX	11	VAL
40	DH	88	GLY
45	DS	71	VAL
46	DU	24	VAL
48	DG	167	VAL
2	AC	56	ILE
2	AC	80	GLY
3	AD	107	GLY
4	AE	56	PRO
4	AE	135	VAL
5	AF	89	VAL
7	AH	33	VAL
9	AJ	8	ILE
16	AS	10	ILE
20	AO	82	ILE
25	BC	150	GLY
25	BC	243	PRO
25	BC	248	GLY
27	BK	93	GLN
28	BP	63	ILE
36	B2	9	VAL
37	BL	65	GLY
45	BS	71	VAL
46	BU	24	VAL
47	BF	88	VAL

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Mol	Chain	Res	Type
48	BG	20	GLY
52	BW	33	GLY
4	CE	133	ILE
7	CH	26	MET
11	CL	3	VAL
6	CG	131	GLY
16	CS	10	ILE
18	CB	28	PRO
25	DC	150	GLY
29	DE	148	ILE
31	D0	46	GLY
37	DL	90	VAL
47	DF	84	ILE
47	DF	88	VAL
49	DR	57	GLY
49	DR	64	VAL
52	DW	21	GLY
4	AE	132	PRO
7	AH	26	MET
8	AI	57	VAL
9	AJ	38	GLY
25	BC	227	VAL
29	BE	64	GLY
35	BV	37	PRO
37	BL	90	VAL
43	BO	49	VAL
49	BR	57	GLY
52	BW	21	GLY
4	CE	135	VAL
10	CK	112	VAL
2	CC	108	PRO
6	CG	18	GLY
25	DC	232	GLY
25	DC	243	PRO
26	DD	194	PRO
29	DE	64	GLY
35	DV	37	PRO
43	DO	28	VAL
48	DG	20	GLY
9	AJ	33	GLY
9	AJ	78	GLU
10	AK	112	VAL

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Mol	Chain	Res	Type
15	AR	38	ILE
18	AB	66	ILE
21	AN	71	GLY
27	BK	103	VAL
29	BE	59	PRO
29	BE	177	PRO
44	BQ	7	VAL
49	BR	64	VAL
49	BR	69	GLY
3	CD	107	GLY
6	CG	92	PRO
25	DC	227	VAL
27	DK	103	VAL
29	DE	59	PRO
43	DO	49	VAL
43	DO	101	GLY
44	DQ	7	VAL
46	DU	10	VAL
4	AE	108	GLY
9	AJ	41	PRO
15	AR	25	ILE
25	BC	7	PRO
26	BD	194	PRO
28	BP	4	ILE
43	BO	28	VAL
47	BF	137	PHE
47	BF	140	ILE
15	CR	25	ILE
15	CR	38	ILE
18	CB	98	GLY
25	DC	7	PRO
28	DP	4	ILE
41	DJ	8	PRO
47	DF	137	PHE
47	DF	140	ILE

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	AC	170/189 (90%)	139 (82%)	31 (18%)	1	11
2	CC	170/189 (90%)	134 (79%)	36 (21%)	1	6
3	AD	172/172 (100%)	145 (84%)	27 (16%)	2	16
3	CD	172/172 (100%)	146 (85%)	26 (15%)	3	17
4	AE	113/125 (90%)	93 (82%)	20 (18%)	2	12
4	CE	113/125 (90%)	93 (82%)	20 (18%)	2	12
5	AF	87/116 (75%)	76 (87%)	11 (13%)	4	22
5	CF	87/116 (75%)	75 (86%)	12 (14%)	3	21
6	AG	123/146 (84%)	102 (83%)	21 (17%)	2	13
6	CG	125/146 (86%)	98 (78%)	27 (22%)	1	6
7	AH	104/104 (100%)	95 (91%)	9 (9%)	10	35
7	CH	104/104 (100%)	95 (91%)	9 (9%)	10	35
8	AI	105/106 (99%)	78 (74%)	27 (26%)	0	4
8	CI	105/106 (99%)	79 (75%)	26 (25%)	0	4
9	AJ	86/90 (96%)	73 (85%)	13 (15%)	3	17
9	CJ	86/90 (96%)	72 (84%)	14 (16%)	2	15
10	AK	90/98 (92%)	73 (81%)	17 (19%)	1	10
10	CK	90/98 (92%)	73 (81%)	17 (19%)	1	10
11	AL	103/103 (100%)	87 (84%)	16 (16%)	2	17
11	CL	103/103 (100%)	87 (84%)	16 (16%)	2	17
12	AM	92/95 (97%)	72 (78%)	20 (22%)	1	6
12	CM	91/95 (96%)	75 (82%)	16 (18%)	2	12
13	AP	65/65 (100%)	61 (94%)	4 (6%)	18	46
13	CP	65/65 (100%)	61 (94%)	4 (6%)	18	46
14	AQ	74/77 (96%)	63 (85%)	11 (15%)	3	18
14	CQ	75/77 (97%)	63 (84%)	12 (16%)	2	16
15	AR	48/64 (75%)	41 (85%)	7 (15%)	3	18
15	CR	48/64 (75%)	40 (83%)	8 (17%)	2	14
16	AS	70/78 (90%)	49 (70%)	21 (30%)	0	2
16	CS	71/78 (91%)	51 (72%)	20 (28%)	0	2
17	AT	65/65 (100%)	51 (78%)	14 (22%)	1	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	CT	65/65 (100%)	51 (78%)	14 (22%)	1	6
18	AB	180/198 (91%)	141 (78%)	39 (22%)	1	6
18	CB	180/198 (91%)	133 (74%)	47 (26%)	0	4
19	AU	44/60 (73%)	31 (70%)	13 (30%)	0	2
19	CU	44/60 (73%)	32 (73%)	12 (27%)	0	3
20	AO	76/77 (99%)	65 (86%)	11 (14%)	3	19
20	CO	76/77 (99%)	61 (80%)	15 (20%)	1	8
21	AN	79/83 (95%)	63 (80%)	16 (20%)	1	7
21	CN	79/83 (95%)	64 (81%)	15 (19%)	1	9
24	BI	109/109 (100%)	108 (99%)	1 (1%)	78	87
24	DI	109/109 (100%)	103 (94%)	6 (6%)	21	50
25	BC	216/217 (100%)	180 (83%)	36 (17%)	2	14
25	DC	216/217 (100%)	181 (84%)	35 (16%)	2	15
26	BD	164/164 (100%)	134 (82%)	30 (18%)	1	11
26	DD	164/164 (100%)	133 (81%)	31 (19%)	1	10
27	BK	102/104 (98%)	76 (74%)	26 (26%)	0	4
27	DK	102/104 (98%)	78 (76%)	24 (24%)	1	5
28	BP	99/99 (100%)	77 (78%)	22 (22%)	1	6
28	DP	99/99 (100%)	77 (78%)	22 (22%)	1	6
29	BE	165/165 (100%)	143 (87%)	22 (13%)	4	22
29	DE	165/165 (100%)	143 (87%)	22 (13%)	4	22
30	BY	48/48 (100%)	39 (81%)	9 (19%)	1	10
30	DY	48/48 (100%)	39 (81%)	9 (19%)	1	10
31	B0	47/47 (100%)	36 (77%)	11 (23%)	1	5
31	D0	47/47 (100%)	37 (79%)	10 (21%)	1	6
32	B4	34/34 (100%)	30 (88%)	4 (12%)	5	24
32	D4	34/34 (100%)	31 (91%)	3 (9%)	10	35
33	B1	45/48 (94%)	39 (87%)	6 (13%)	4	22
33	D1	45/48 (94%)	39 (87%)	6 (13%)	4	22
34	B3	51/51 (100%)	45 (88%)	6 (12%)	5	24
34	D3	51/51 (100%)	45 (88%)	6 (12%)	5	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
35	BV	78/78 (100%)	64 (82%)	14 (18%)	2	12
35	DV	78/78 (100%)	64 (82%)	14 (18%)	2	12
36	B2	38/38 (100%)	32 (84%)	6 (16%)	2	16
36	D2	38/38 (100%)	33 (87%)	5 (13%)	4	22
37	BL	102/103 (99%)	89 (87%)	13 (13%)	4	22
37	DL	102/103 (99%)	90 (88%)	12 (12%)	5	24
38	BM	109/109 (100%)	86 (79%)	23 (21%)	1	6
38	DM	109/109 (100%)	86 (79%)	23 (21%)	1	6
39	BX	55/55 (100%)	43 (78%)	12 (22%)	1	6
39	DX	55/55 (100%)	45 (82%)	10 (18%)	1	11
40	BH	114/114 (100%)	81 (71%)	33 (29%)	0	2
40	DH	114/114 (100%)	89 (78%)	25 (22%)	1	6
41	BJ	116/116 (100%)	95 (82%)	21 (18%)	1	11
41	DJ	116/116 (100%)	96 (83%)	20 (17%)	2	13
42	BN	100/103 (97%)	87 (87%)	13 (13%)	4	22
42	DN	100/103 (97%)	87 (87%)	13 (13%)	4	22
43	BO	86/87 (99%)	72 (84%)	14 (16%)	2	15
43	DO	86/87 (99%)	72 (84%)	14 (16%)	2	15
44	BQ	89/89 (100%)	74 (83%)	15 (17%)	2	14
44	DQ	89/89 (100%)	74 (83%)	15 (17%)	2	14
45	BS	93/93 (100%)	80 (86%)	13 (14%)	3	20
45	DS	93/93 (100%)	80 (86%)	13 (14%)	3	20
46	BU	83/84 (99%)	69 (83%)	14 (17%)	2	14
46	DU	83/84 (99%)	69 (83%)	14 (17%)	2	14
47	BF	149/149 (100%)	117 (78%)	32 (22%)	1	6
47	DF	149/149 (100%)	116 (78%)	33 (22%)	1	6
48	BG	137/137 (100%)	112 (82%)	25 (18%)	1	11
48	DG	137/137 (100%)	112 (82%)	25 (18%)	1	11
49	BR	84/84 (100%)	71 (84%)	13 (16%)	2	17
49	DR	84/84 (100%)	73 (87%)	11 (13%)	4	22
50	BT	80/84 (95%)	59 (74%)	21 (26%)	0	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
50	DT	80/84 (95%)	60 (75%)	20 (25%)	0	4
51	BZ	67/68 (98%)	53 (79%)	14 (21%)	1	7
51	DZ	67/68 (98%)	53 (79%)	14 (21%)	1	7
52	BW	59/62 (95%)	42 (71%)	17 (29%)	0	2
52	DW	59/62 (95%)	42 (71%)	17 (29%)	0	2
All	All	9333/9700 (96%)	7661 (82%)	1672 (18%)	2	12

All (1672) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	AC	2	GLN
2	AC	13	ILE
2	AC	15	LYS
2	AC	17	TRP
2	AC	19	SER
2	AC	27	GLU
2	AC	40	GLN
2	AC	41	TYR
2	AC	48	LYS
2	AC	62	SER
2	AC	66	THR
2	AC	69	THR
2	AC	78	LYS
2	AC	79	LYS
2	AC	87	ARG
2	AC	96	VAL
2	AC	101	ASN
2	AC	105	VAL
2	AC	110	LEU
2	AC	113	LYS
2	AC	118	SER
2	AC	126	ARG
2	AC	128	MET
2	AC	139	ASN
2	AC	152	VAL
2	AC	155	ARG
2	AC	168	ARG
2	AC	171	ARG
2	AC	172	VAL
2	AC	183	TYR

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Mol	Chain	Res	Type
2	AC	184	ASN
3	AD	2	ARG
3	AD	12	ARG
3	AD	18	LEU
3	AD	25	ARG
3	AD	27	ILE
3	AD	28	ASP
3	AD	34	GLU
3	AD	46	ARG
3	AD	49	ASP
3	AD	55	ARG
3	AD	87	GLU
3	AD	104	MET
3	AD	106	PHE
3	AD	114	ARG
3	AD	118	SER
3	AD	146	GLU
3	AD	147	LYS
3	AD	158	LEU
3	AD	160	LEU
3	AD	176	LYS
3	AD	189	ASP
3	AD	190	LEU
3	AD	196	GLU
3	AD	197	HIS
3	AD	198	LEU
3	AD	199	ILE
3	AD	200	VAL
4	AE	14	LEU
4	AE	23	THR
4	AE	30	PHE
4	AE	45	VAL
4	AE	51	LYS
4	AE	60	GLN
4	AE	61	LYS
4	AE	63	MET
4	AE	68	ARG
4	AE	72	ASN
4	AE	77	ASN
4	AE	81	GLN
4	AE	92	ARG
4	AE	115	GLU

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Mol	Chain	Res	Type
4	AE	123	LEU
4	AE	125	LYS
4	AE	127	TYR
4	AE	144	GLU
4	AE	147	ASN
4	AE	156	ARG
5	AF	6	ILE
5	AF	16	GLU
5	AF	54	LEU
5	AF	56	LYS
5	AF	62	MET
5	AF	69	GLU
5	AF	73	GLU
5	AF	86	ARG
5	AF	96	VAL
5	AF	97	THR
5	AF	100	SER
6	AG	5	VAL
6	AG	6	ILE
6	AG	8	GLN
6	AG	11	ILE
6	AG	12	LEU
6	AG	16	LYS
6	AG	21	LEU
6	AG	29	LEU
6	AG	55	LYS
6	AG	57	GLU
6	AG	58	LEU
6	AG	59	GLU
6	AG	68	VAL
6	AG	72	VAL
6	AG	89	GLU
6	AG	94	ARG
6	AG	96	ASN
6	AG	117	LEU
6	AG	124	SER
6	AG	138	GLU
6	AG	143	MET
7	AH	24	VAL
7	AH	53	ASP
7	AH	54	THR
7	AH	55	LYS

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Mol	Chain	Res	Type
7	AH	72	GLU
7	AH	82	LEU
7	AH	100	ILE
7	AH	105	THR
7	AH	113	ARG
8	AI	12	LYS
8	AI	13	SER
8	AI	14	SER
8	AI	26	LYS
8	AI	30	ASN
8	AI	31	GLN
8	AI	36	GLN
8	AI	37	TYR
8	AI	42	THR
8	AI	45	MET
8	AI	46	VAL
8	AI	62	LEU
8	AI	67	LYS
8	AI	87	MET
8	AI	89	TYR
8	AI	90	ASP
8	AI	91	GLU
8	AI	94	ARG
8	AI	99	LYS
8	AI	105	ARG
8	AI	106	ASP
8	AI	109	GLN
8	AI	112	ARG
8	AI	122	ARG
8	AI	123	ARG
8	AI	126	PHE
8	AI	128	LYS
9	AJ	15	HIS
9	AJ	17	LEU
9	AJ	19	ASP
9	AJ	32	THR
9	AJ	60	ASP
9	AJ	64	GLN
9	AJ	75	ASP
9	AJ	80	THR
9	AJ	81	GLU
9	AJ	83	THR

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Mol	Chain	Res	Type
9	AJ	88	MET
9	AJ	92	LEU
9	AJ	102	LEU
10	AK	26	PHE
10	AK	28	ASN
10	AK	34	THR
10	AK	35	ASP
10	AK	52	ARG
10	AK	69	CYS
10	AK	75	GLU
10	AK	80	ASN
10	AK	84	MET
10	AK	100	ASN
10	AK	105	ARG
10	AK	106	ILE
10	AK	110	THR
10	AK	115	ILE
10	AK	117	HIS
10	AK	118	ASN
10	AK	127	ARG
11	AL	9	LYS
11	AL	14	LYS
11	AL	17	LYS
11	AL	28	GLN
11	AL	35	ARG
11	AL	43	LYS
11	AL	48	LEU
11	AL	49	ARG
11	AL	86	VAL
11	AL	87	LYS
11	AL	93	ARG
11	AL	95	HIS
11	AL	96	THR
11	AL	107	LYS
11	AL	108	ASP
11	AL	122	LYS
12	AM	2	ARG
12	AM	16	ILE
12	AM	26	LYS
12	AM	28	ARG
12	AM	44	ILE
12	AM	45	SER

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Mol	Chain	Res	Type
12	AM	46	GLU
12	AM	47	LEU
12	AM	57	ASP
12	AM	62	PHE
12	AM	67	ASP
12	AM	71	GLU
12	AM	77	LYS
12	AM	89	ARG
12	AM	91	ARG
12	AM	92	ARG
12	AM	96	VAL
12	AM	99	GLN
12	AM	101	THR
12	AM	102	LYS
13	AP	8	ARG
13	AP	28	ARG
13	AP	45	GLU
13	AP	51	ARG
14	AQ	5	ARG
14	AQ	6	THR
14	AQ	15	LYS
14	AQ	27	PHE
14	AQ	39	ARG
14	AQ	48	GLU
14	AQ	60	ILE
14	AQ	66	LEU
14	AQ	74	LEU
14	AQ	79	GLU
14	AQ	80	LYS
15	AR	23	LYS
15	AR	30	ASN
15	AR	38	ILE
15	AR	42	ARG
15	AR	46	THR
15	AR	71	ASP
15	AR	72	ARG
16	AS	2	ARG
16	AS	4	LEU
16	AS	5	LYS
16	AS	10	ILE
16	AS	11	ASP
16	AS	14	LEU

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Mol	Chain	Res	Type
16	AS	20	LYS
16	AS	26	ASP
16	AS	27	LYS
16	AS	28	LYS
16	AS	33	TRP
16	AS	34	SER
16	AS	36	ARG
16	AS	38	THR
16	AS	40	PHE
16	AS	47	THR
16	AS	54	ARG
16	AS	59	VAL
16	AS	64	GLU
16	AS	70	LEU
16	AS	76	THR
17	AT	2	ASN
17	AT	4	LYS
17	AT	13	SER
17	AT	14	GLU
17	AT	35	TYR
17	AT	43	LYS
17	AT	51	ASN
17	AT	52	GLU
17	AT	53	MET
17	AT	59	ARG
17	AT	67	HIS
17	AT	69	ASN
17	AT	73	ARG
17	AT	79	THR
18	AB	8	MET
18	AB	14	HIS
18	AB	18	GLN
18	AB	23	ASN
18	AB	31	PHE
18	AB	38	HIS
18	AB	41	ASN
18	AB	42	LEU
18	AB	43	GLU
18	AB	46	VAL
18	AB	48	MET
18	AB	61	SER
18	AB	62	ARG

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Mol	Chain	Res	Type
18	AB	67	LEU
18	AB	68	PHE
18	AB	73	ARG
18	AB	80	LYS
18	AB	81	ASP
18	AB	84	LEU
18	AB	88	GLN
18	AB	94	ARG
18	AB	95	TRP
18	AB	100	LEU
18	AB	104	LYS
18	AB	107	ARG
18	AB	125	PHE
18	AB	127	LYS
18	AB	131	LYS
18	AB	146	SER
18	AB	160	LEU
18	AB	169	HIS
18	AB	183	PHE
18	AB	187	ASP
18	AB	188	THR
18	AB	196	ASP
18	AB	202	ASN
18	AB	203	ASP
18	AB	213	LEU
18	AB	221	ARG
19	AU	7	GLU
19	AU	11	PHE
19	AU	15	LEU
19	AU	16	ARG
19	AU	18	PHE
19	AU	19	LYS
19	AU	20	ARG
19	AU	22	CYS
19	AU	24	LYS
19	AU	33	ARG
19	AU	34	ARG
19	AU	36	PHE
19	AU	44	ARG
20	AO	15	PHE
20	AO	17	ARG
20	AO	20	ASN

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Mol	Chain	Res	Type
20	AO	32	LEU
20	AO	33	THR
20	AO	40	GLN
20	AO	42	HIS
20	AO	66	LEU
20	AO	70	LEU
20	AO	77	ARG
20	AO	88	ARG
21	AN	19	TYR
21	AN	20	PHE
21	AN	30	ILE
21	AN	44	VAL
21	AN	47	LEU
21	AN	48	GLN
21	AN	49	THR
21	AN	50	LEU
21	AN	52	ARG
21	AN	53	ASP
21	AN	59	GLN
21	AN	65	GLN
21	AN	74	ARG
21	AN	80	ARG
21	AN	84	ARG
21	AN	100	TRP
24	BI	96	LYS
25	BC	5	CYS
25	BC	9	SER
25	BC	12	ARG
25	BC	35	LYS
25	BC	43	ASN
25	BC	45	ASN
25	BC	62	ARG
25	BC	63	ILE
25	BC	65	ASP
25	BC	66	PHE
25	BC	69	ASN
25	BC	71	ASP
25	BC	82	TYR
25	BC	86	ARG
25	BC	100	ARG
25	BC	113	ASP
25	BC	123	ILE

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Mol	Chain	Res	Type
25	BC	134	ILE
25	BC	152	GLN
25	BC	155	ARG
25	BC	166	ARG
25	BC	173	LEU
25	BC	176	ARG
25	BC	180	MET
25	BC	181	ARG
25	BC	186	ASP
25	BC	188	ARG
25	BC	191	LEU
25	BC	201	LEU
25	BC	224	MET
25	BC	239	PHE
25	BC	245	THR
25	BC	249	VAL
25	BC	251	THR
25	BC	255	LYS
25	BC	269	ARG
26	BD	1	MET
26	BD	11	MET
26	BD	13	ARG
26	BD	17	GLU
26	BD	33	ARG
26	BD	36	GLN
26	BD	40	LEU
26	BD	46	ARG
26	BD	52	THR
26	BD	55	LYS
26	BD	58	ASN
26	BD	62	LYS
26	BD	81	GLU
26	BD	82	PHE
26	BD	84	LEU
26	BD	91	THR
26	BD	98	VAL
26	BD	106	LYS
26	BD	108	ASP
26	BD	118	PHE
26	BD	128	ARG
26	BD	131	ASP
26	BD	138	LEU

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Mol	Chain	Res	Type
26	BD	141	ARG
26	BD	148	GLN
26	BD	154	LYS
26	BD	159	LYS
26	BD	183	GLU
26	BD	201	LEU
26	BD	204	LYS
27	BK	6	THR
27	BK	18	ARG
27	BK	19	VAL
27	BK	21	CYS
27	BK	25	LEU
27	BK	32	TYR
27	BK	37	ASP
27	BK	38	ILE
27	BK	47	ILE
27	BK	52	VAL
27	BK	54	LYS
27	BK	65	THR
27	BK	66	LYS
27	BK	72	PRO
27	BK	73	ASP
27	BK	79	PHE
27	BK	86	LEU
27	BK	88	ASN
27	BK	97	THR
27	BK	104	THR
27	BK	105	ARG
27	BK	107	LEU
27	BK	108	ARG
27	BK	111	LYS
27	BK	113	MET
27	BK	120	PRO
28	BP	3	ILE
28	BP	6	GLN
28	BP	19	PHE
28	BP	20	ARG
28	BP	25	VAL
28	BP	33	GLU
28	BP	43	GLU
28	BP	52	ARG
28	BP	58	PHE

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Mol	Chain	Res	Type
28	BP	59	THR
28	BP	61	ARG
28	BP	65	ASN
28	BP	80	VAL
28	BP	83	ILE
28	BP	84	SER
28	BP	95	LYS
28	BP	99	LEU
28	BP	100	ARG
28	BP	101	GLU
28	BP	111	GLU
28	BP	112	ARG
28	BP	114	ASN
29	BE	5	LEU
29	BE	6	LYS
29	BE	7	ASP
29	BE	9	GLN
29	BE	13	THR
29	BE	16	GLU
29	BE	17	THR
29	BE	24	ASN
29	BE	58	LYS
29	BE	61	ARG
29	BE	62	GLN
29	BE	63	LYS
29	BE	78	TRP
29	BE	108	ILE
29	BE	112	LEU
29	BE	118	LEU
29	BE	122	GLU
29	BE	123	LYS
29	BE	144	GLU
29	BE	145	ASP
29	BE	147	LEU
29	BE	198	GLU
30	BY	2	LYS
30	BY	6	ILE
30	BY	10	ARG
30	BY	15	ARG
30	BY	23	LEU
30	BY	24	LEU
30	BY	37	ARG

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Mol	Chain	Res	Type
30	BY	43	ILE
30	BY	53	MET
31	B0	5	ASN
31	B0	26	SER
31	B0	27	LEU
31	B0	28	SER
31	B0	35	GLU
31	B0	37	HIS
31	B0	38	LEU
31	B0	41	HIS
31	B0	45	ASP
31	B0	47	TYR
31	B0	51	ARG
32	B4	3	VAL
32	B4	9	LYS
32	B4	24	ARG
32	B4	35	GLN
33	B1	6	GLU
33	B1	8	ILE
33	B1	9	LYS
33	B1	16	THR
33	B1	31	GLU
33	B1	35	LEU
34	B3	7	ARG
34	B3	14	LYS
34	B3	18	LYS
34	B3	56	LEU
34	B3	57	VAL
34	B3	58	ILE
35	BV	11	GLU
35	BV	18	ARG
35	BV	24	ASN
35	BV	34	LYS
35	BV	40	ILE
35	BV	42	LEU
35	BV	51	GLN
35	BV	53	LYS
35	BV	68	LYS
35	BV	70	ILE
35	BV	75	GLN
35	BV	82	TYR
35	BV	87	GLN

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Mol	Chain	Res	Type
35	BV	90	ASP
36	B2	3	ARG
36	B2	4	THR
36	B2	22	MET
36	B2	34	ARG
36	B2	39	ARG
36	B2	46	LYS
37	BL	4	ASN
37	BL	19	LEU
37	BL	21	ARG
37	BL	46	VAL
37	BL	64	PHE
37	BL	69	ARG
37	BL	91	ASP
37	BL	92	LEU
37	BL	94	THR
37	BL	95	LEU
37	BL	118	THR
37	BL	122	VAL
37	BL	123	ARG
38	BM	1	MET
38	BM	20	LEU
38	BM	25	ASP
38	BM	42	THR
38	BM	46	ILE
38	BM	50	ARG
38	BM	55	ARG
38	BM	59	ARG
38	BM	60	GLN
38	BM	63	ILE
38	BM	65	ILE
38	BM	70	ASP
38	BM	75	GLU
38	BM	78	LEU
38	BM	81	ARG
38	BM	82	MET
38	BM	88	ASN
38	BM	90	GLU
38	BM	93	VAL
38	BM	110	GLU
38	BM	111	GLU
38	BM	115	GLU

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Mol	Chain	Res	Type
38	BM	131	VAL
39	BX	2	LYS
39	BX	7	ARG
39	BX	16	THR
39	BX	17	GLU
39	BX	20	ASN
39	BX	21	LEU
39	BX	22	LEU
39	BX	25	GLN
39	BX	37	LEU
39	BX	57	LEU
39	BX	58	ASN
39	BX	59	GLU
40	BH	1	MET
40	BH	3	VAL
40	BH	12	LEU
40	BH	15	LEU
40	BH	17	ASP
40	BH	25	TYR
40	BH	28	ASN
40	BH	43	ASN
40	BH	44	ILE
40	BH	50	ARG
40	BH	54	LEU
40	BH	55	GLU
40	BH	57	LYS
40	BH	62	LEU
40	BH	66	ASN
40	BH	68	ARG
40	BH	70	GLU
40	BH	71	LYS
40	BH	73	ASN
40	BH	87	GLU
40	BH	89	LYS
40	BH	90	LEU
40	BH	98	ASP
40	BH	103	VAL
40	BH	104	THR
40	BH	109	GLU
40	BH	110	VAL
40	BH	112	LYS
40	BH	116	ARG

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Mol	Chain	Res	Type
40	BH	128	HIS
40	BH	130	VAL
40	BH	135	HIS
40	BH	146	VAL
41	BJ	2	LYS
41	BJ	3	THR
41	BJ	5	THR
41	BJ	12	LYS
41	BJ	21	THR
41	BJ	31	GLU
41	BJ	43	GLU
41	BJ	44	TYR
41	BJ	50	THR
41	BJ	54	ILE
41	BJ	65	THR
41	BJ	72	LYS
41	BJ	73	VAL
41	BJ	95	ARG
41	BJ	106	LYS
41	BJ	108	MET
41	BJ	124	VAL
41	BJ	129	GLU
41	BJ	131	ASN
41	BJ	136	GLN
41	BJ	140	LEU
42	BN	2	ARG
42	BN	3	HIS
42	BN	11	ASN
42	BN	22	ARG
42	BN	33	ILE
42	BN	35	LYS
42	BN	46	ARG
42	BN	51	LEU
42	BN	57	THR
42	BN	69	ARG
42	BN	100	CYS
42	BN	114	GLU
42	BN	120	GLU
43	BO	3	LYS
43	BO	9	ARG
43	BO	21	LEU
43	BO	31	THR

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Mol	Chain	Res	Type
43	BO	33	ARG
43	BO	35	ILE
43	BO	43	ASN
43	BO	46	GLU
43	BO	78	VAL
43	BO	84	GLU
43	BO	98	GLN
43	BO	104	GLN
43	BO	106	LEU
43	BO	116	GLN
44	BQ	5	ARG
44	BQ	10	ARG
44	BQ	13	HIS
44	BQ	14	LYS
44	BQ	19	GLN
44	BQ	29	ARG
44	BQ	48	ASP
44	BQ	59	LEU
44	BQ	69	ARG
44	BQ	79	ILE
44	BQ	83	LYS
44	BQ	88	GLU
44	BQ	90	ASP
44	BQ	96	ASP
44	BQ	105	PHE
45	BS	1	MET
45	BS	6	LYS
45	BS	17	VAL
45	BS	18	ARG
45	BS	22	ASP
45	BS	36	LEU
45	BS	40	ASN
45	BS	66	ILE
45	BS	72	THR
45	BS	84	ARG
45	BS	86	MET
45	BS	88	ARG
45	BS	101	SER
46	BU	7	ASP
46	BU	11	ILE
46	BU	18	LYS
46	BU	26	ASN

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Mol	Chain	Res	Type
46	BU	35	VAL
46	BU	45	GLN
46	BU	46	LYS
46	BU	51	LEU
46	BU	60	LYS
46	BU	73	ASN
46	BU	78	LYS
46	BU	85	ARG
46	BU	87	GLU
46	BU	102	ILE
47	BF	2	LYS
47	BF	9	ASP
47	BF	18	GLU
47	BF	29	ARG
47	BF	35	LEU
47	BF	49	LEU
47	BF	50	ASP
47	BF	56	LEU
47	BF	62	GLN
47	BF	66	ILE
47	BF	68	LYS
47	BF	76	PHE
47	BF	91	ARG
47	BF	93	GLU
47	BF	94	ARG
47	BF	97	GLU
47	BF	100	GLU
47	BF	111	ARG
47	BF	121	PHE
47	BF	129	MET
47	BF	134	GLN
47	BF	137	PHE
47	BF	138	PRO
47	BF	141	ASP
47	BF	142	TYR
47	BF	143	ASP
47	BF	146	ASP
47	BF	147	ARG
47	BF	149	ARG
47	BF	156	THR
47	BF	177	ARG
47	BF	178	LYS

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Mol	Chain	Res	Type
48	BG	14	VAL
48	BG	17	LYS
48	BG	19	ASN
48	BG	26	LYS
48	BG	29	ASN
48	BG	31	GLU
48	BG	34	ARG
48	BG	36	LEU
48	BG	46	ASP
48	BG	50	THR
48	BG	54	ARG
48	BG	68	ARG
48	BG	84	LYS
48	BG	85	LYS
48	BG	86	LEU
48	BG	94	ARG
48	BG	120	ILE
48	BG	123	GLU
48	BG	138	GLN
48	BG	148	ARG
48	BG	162	ARG
48	BG	166	GLU
48	BG	167	VAL
48	BG	169	ARG
48	BG	176	LYS
49	BR	5	PHE
49	BR	12	HIS
49	BR	19	THR
49	BR	22	LEU
49	BR	27	ILE
49	BR	39	LEU
49	BR	41	ILE
49	BR	48	LYS
49	BR	53	PHE
49	BR	66	HIS
49	BR	70	GLU
49	BR	72	VAL
49	BR	99	THR
50	BT	2	ILE
50	BT	3	ARG
50	BT	4	GLU
50	BT	6	ARG

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Mol	Chain	Res	Type
50	BT	9	LYS
50	BT	11	LEU
50	BT	12	ARG
50	BT	22	THR
50	BT	29	THR
50	BT	32	LEU
50	BT	50	LEU
50	BT	54	GLU
50	BT	61	LEU
50	BT	64	LYS
50	BT	68	LYS
50	BT	69	ARG
50	BT	73	ARG
50	BT	76	ARG
50	BT	79	ASP
50	BT	87	LEU
50	BT	92	ASN
51	BZ	4	VAL
51	BZ	13	VAL
51	BZ	27	ARG
51	BZ	28	ARG
51	BZ	29	PHE
51	BZ	30	LEU
51	BZ	33	LEU
51	BZ	37	ARG
51	BZ	46	PHE
51	BZ	49	LEU
51	BZ	60	ASP
51	BZ	65	ASP
51	BZ	77	LYS
51	BZ	78	TYR
52	BW	10	ARG
52	BW	16	GLU
52	BW	18	LYS
52	BW	19	ARG
52	BW	23	LYS
52	BW	24	ARG
52	BW	25	PHE
52	BW	31	LEU
52	BW	38	ARG
52	BW	39	GLN
52	BW	40	ARG

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Mol	Chain	Res	Type
52	BW	44	PHE
52	BW	50	VAL
52	BW	61	LYS
52	BW	75	ASN
52	BW	76	ARG
52	BW	77	LYS
3	CD	2	ARG
3	CD	12	ARG
3	CD	18	LEU
3	CD	25	ARG
3	CD	27	ILE
3	CD	28	ASP
3	CD	34	GLU
3	CD	46	ARG
3	CD	49	ASP
3	CD	55	ARG
3	CD	87	GLU
3	CD	104	MET
3	CD	106	PHE
3	CD	114	ARG
3	CD	146	GLU
3	CD	147	LYS
3	CD	158	LEU
3	CD	160	LEU
3	CD	176	LYS
3	CD	189	ASP
3	CD	190	LEU
3	CD	196	GLU
3	CD	197	HIS
3	CD	198	LEU
3	CD	199	ILE
3	CD	200	VAL
4	CE	14	LEU
4	CE	23	THR
4	CE	30	PHE
4	CE	45	VAL
4	CE	51	LYS
4	CE	60	GLN
4	CE	61	LYS
4	CE	63	MET
4	CE	68	ARG
4	CE	72	ASN

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Mol	Chain	Res	Type
4	CE	77	ASN
4	CE	81	GLN
4	CE	92	ARG
4	CE	115	GLU
4	CE	123	LEU
4	CE	125	LYS
4	CE	127	TYR
4	CE	144	GLU
4	CE	147	ASN
4	CE	156	ARG
5	CF	6	ILE
5	CF	16	GLU
5	CF	54	LEU
5	CF	56	LYS
5	CF	62	MET
5	CF	69	GLU
5	CF	72	ASP
5	CF	73	GLU
5	CF	86	ARG
5	CF	96	VAL
5	CF	97	THR
5	CF	100	SER
7	CH	24	VAL
7	CH	25	THR
7	CH	53	ASP
7	CH	54	THR
7	CH	55	LYS
7	CH	72	GLU
7	CH	82	LEU
7	CH	100	ILE
7	CH	113	ARG
10	CK	26	PHE
10	CK	28	ASN
10	CK	34	THR
10	CK	35	ASP
10	CK	52	ARG
10	CK	69	CYS
10	CK	75	GLU
10	CK	80	ASN
10	CK	84	MET
10	CK	100	ASN
10	CK	105	ARG

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Mol	Chain	Res	Type
10	CK	106	ILE
10	CK	110	THR
10	CK	115	ILE
10	CK	117	HIS
10	CK	118	ASN
10	CK	127	ARG
11	CL	9	LYS
11	CL	14	LYS
11	CL	17	LYS
11	CL	28	GLN
11	CL	35	ARG
11	CL	43	LYS
11	CL	48	LEU
11	CL	49	ARG
11	CL	86	VAL
11	CL	87	LYS
11	CL	93	ARG
11	CL	95	HIS
11	CL	96	THR
11	CL	107	LYS
11	CL	108	ASP
11	CL	122	LYS
13	CP	8	ARG
13	CP	28	ARG
13	CP	45	GLU
13	CP	51	ARG
14	CQ	5	ARG
14	CQ	6	THR
14	CQ	8	GLN
14	CQ	15	LYS
14	CQ	27	PHE
14	CQ	39	ARG
14	CQ	48	GLU
14	CQ	60	ILE
14	CQ	66	LEU
14	CQ	74	LEU
14	CQ	79	GLU
14	CQ	80	LYS
15	CR	23	LYS
15	CR	30	ASN
15	CR	38	ILE
15	CR	42	ARG

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Mol	Chain	Res	Type
15	CR	46	THR
15	CR	71	ASP
15	CR	72	ARG
15	CR	73	HIS
17	CT	2	ASN
17	CT	4	LYS
17	CT	13	SER
17	CT	14	GLU
17	CT	35	TYR
17	CT	43	LYS
17	CT	51	ASN
17	CT	52	GLU
17	CT	53	MET
17	CT	59	ARG
17	CT	67	HIS
17	CT	69	ASN
17	CT	73	ARG
17	CT	79	THR
19	CU	11	PHE
19	CU	15	LEU
19	CU	16	ARG
19	CU	18	PHE
19	CU	19	LYS
19	CU	20	ARG
19	CU	22	CYS
19	CU	24	LYS
19	CU	33	ARG
19	CU	34	ARG
19	CU	36	PHE
19	CU	44	ARG
2	CC	2	GLN
2	CC	10	ARG
2	CC	13	ILE
2	CC	17	TRP
2	CC	20	THR
2	CC	27	GLU
2	CC	31	ASN
2	CC	39	ARG
2	CC	44	LYS
2	CC	48	LYS
2	CC	50	SER
2	CC	56	ILE

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Mol	Chain	Res	Type
2	CC	61	LYS
2	CC	63	ILE
2	CC	78	LYS
2	CC	86	LEU
2	CC	88	LYS
2	CC	104	GLU
2	CC	106	ARG
2	CC	107	LYS
2	CC	109	GLU
2	CC	117	ASP
2	CC	120	THR
2	CC	126	ARG
2	CC	128	MET
2	CC	130	ARG
2	CC	138	GLN
2	CC	143	LEU
2	CC	156	LEU
2	CC	165	GLU
2	CC	168	ARG
2	CC	175	HIS
2	CC	186	SER
2	CC	192	TYR
2	CC	201	ILE
2	CC	205	GLU
6	CG	2	ARG
6	CG	11	ILE
6	CG	12	LEU
6	CG	21	LEU
6	CG	26	VAL
6	CG	47	GLU
6	CG	55	LYS
6	CG	56	SER
6	CG	59	GLU
6	CG	62	GLU
6	CG	67	ASN
6	CG	75	LYS
6	CG	78	ARG
6	CG	91	ARG
6	CG	93	VAL
6	CG	94	ARG
6	CG	96	ASN
6	CG	105	GLU

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Mol	Chain	Res	Type
6	CG	108	ARG
6	CG	109	LYS
6	CG	112	ASP
6	CG	115	MET
6	CG	124	SER
6	CG	128	GLU
6	CG	134	VAL
6	CG	137	ARG
6	CG	152	HIS
8	CI	3	ASN
8	CI	11	ARG
8	CI	13	SER
8	CI	27	ILE
8	CI	30	ASN
8	CI	35	GLU
8	CI	42	THR
8	CI	44	ARG
8	CI	45	MET
8	CI	46	VAL
8	CI	56	MET
8	CI	58	GLU
8	CI	59	LYS
8	CI	60	LEU
8	CI	62	LEU
8	CI	65	THR
8	CI	67	LYS
8	CI	74	GLN
8	CI	80	HIS
8	CI	93	LEU
8	CI	96	GLU
8	CI	108	ARG
8	CI	109	GLN
8	CI	114	LYS
8	CI	123	ARG
8	CI	129	ARG
9	CJ	24	GLU
9	CJ	37	ARG
9	CJ	47	GLU
9	CJ	64	GLN
9	CJ	72	ARG
9	CJ	73	LEU
9	CJ	78	GLU

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Mol	Chain	Res	Type
9	CJ	80	THR
9	CJ	81	GLU
9	CJ	82	LYS
9	CJ	89	ARG
9	CJ	90	LEU
9	CJ	92	LEU
9	CJ	99	GLN
12	CM	2	ARG
12	CM	8	ILE
12	CM	22	TYR
12	CM	38	ILE
12	CM	40	GLU
12	CM	41	ASP
12	CM	43	LYS
12	CM	47	LEU
12	CM	54	THR
12	CM	55	LEU
12	CM	57	ASP
12	CM	64	VAL
12	CM	65	GLU
12	CM	91	ARG
12	CM	97	ARG
12	CM	106	ARG
21	CN	15	LEU
21	CN	41	TRP
21	CN	45	LEU
21	CN	47	LEU
21	CN	49	THR
21	CN	50	LEU
21	CN	52	ARG
21	CN	55	SER
21	CN	58	ARG
21	CN	63	CYS
21	CN	64	ARG
21	CN	65	GLN
21	CN	76	PHE
21	CN	81	ILE
21	CN	96	LYS
16	CS	2	ARG
16	CS	4	LEU
16	CS	5	LYS
16	CS	10	ILE

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Mol	Chain	Res	Type
16	CS	12	LEU
16	CS	13	HIS
16	CS	15	LEU
16	CS	23	GLU
16	CS	26	ASP
16	CS	27	LYS
16	CS	28	LYS
16	CS	36	ARG
16	CS	38	THR
16	CS	39	ILE
16	CS	47	THR
16	CS	56	HIS
16	CS	59	VAL
16	CS	60	PHE
16	CS	64	GLU
16	CS	65	MET
18	CB	8	MET
18	CB	9	LEU
18	CB	19	THR
18	CB	22	TRP
18	CB	23	ASN
18	CB	27	LYS
18	CB	35	ASN
18	CB	42	LEU
18	CB	43	GLU
18	CB	44	LYS
18	CB	46	VAL
18	CB	50	ASN
18	CB	57	ASN
18	CB	58	LYS
18	CB	62	ARG
18	CB	72	LYS
18	CB	84	LEU
18	CB	86	CYS
18	CB	92	ASN
18	CB	93	HIS
18	CB	95	TRP
18	CB	103	TRP
18	CB	104	LYS
18	CB	113	LEU
18	CB	121	GLN
18	CB	122	ASP

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Mol	Chain	Res	Type
18	CB	126	ASP
18	CB	127	LYS
18	CB	136	ARG
18	CB	137	THR
18	CB	144	GLU
18	CB	152	ASP
18	CB	160	LEU
18	CB	163	ILE
18	CB	166	ASP
18	CB	169	HIS
18	CB	176	ASN
18	CB	178	LEU
18	CB	187	ASP
18	CB	193	ASP
18	CB	196	ASP
18	CB	198	VAL
18	CB	199	ILE
18	CB	210	THR
18	CB	211	LEU
18	CB	212	TYR
18	CB	221	ARG
20	CO	4	SER
20	CO	6	GLU
20	CO	18	ASP
20	CO	33	THR
20	CO	35	GLN
20	CO	37	ASN
20	CO	53	ARG
20	CO	54	ARG
20	CO	59	MET
20	CO	62	GLN
20	CO	64	ARG
20	CO	66	LEU
20	CO	70	LEU
20	CO	71	LYS
20	CO	88	ARG
24	DI	2	LYS
24	DI	54	ILE
24	DI	91	LYS
24	DI	99	LYS
24	DI	121	ILE
24	DI	140	GLU

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Mol	Chain	Res	Type
25	DC	5	CYS
25	DC	9	SER
25	DC	12	ARG
25	DC	35	LYS
25	DC	43	ASN
25	DC	45	ASN
25	DC	62	ARG
25	DC	63	ILE
25	DC	65	ASP
25	DC	66	PHE
25	DC	69	ASN
25	DC	71	ASP
25	DC	82	TYR
25	DC	86	ARG
25	DC	100	ARG
25	DC	113	ASP
25	DC	123	ILE
25	DC	134	ILE
25	DC	152	GLN
25	DC	155	ARG
25	DC	166	ARG
25	DC	173	LEU
25	DC	176	ARG
25	DC	180	MET
25	DC	181	ARG
25	DC	186	ASP
25	DC	188	ARG
25	DC	191	LEU
25	DC	201	LEU
25	DC	224	MET
25	DC	239	PHE
25	DC	245	THR
25	DC	249	VAL
25	DC	251	THR
25	DC	269	ARG
26	DD	1	MET
26	DD	11	MET
26	DD	13	ARG
26	DD	17	GLU
26	DD	33	ARG
26	DD	36	GLN
26	DD	40	LEU

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Mol	Chain	Res	Type
26	DD	46	ARG
26	DD	55	LYS
26	DD	58	ASN
26	DD	62	LYS
26	DD	81	GLU
26	DD	82	PHE
26	DD	84	LEU
26	DD	91	THR
26	DD	98	VAL
26	DD	106	LYS
26	DD	108	ASP
26	DD	118	PHE
26	DD	128	ARG
26	DD	131	ASP
26	DD	138	LEU
26	DD	141	ARG
26	DD	146	ILE
26	DD	148	GLN
26	DD	154	LYS
26	DD	159	LYS
26	DD	183	GLU
26	DD	199	SER
26	DD	201	LEU
26	DD	204	LYS
27	DK	6	THR
27	DK	18	ARG
27	DK	19	VAL
27	DK	21	CYS
27	DK	25	LEU
27	DK	32	TYR
27	DK	38	ILE
27	DK	47	ILE
27	DK	52	VAL
27	DK	54	LYS
27	DK	65	THR
27	DK	66	LYS
27	DK	72	PRO
27	DK	73	ASP
27	DK	79	PHE
27	DK	86	LEU
27	DK	88	ASN
27	DK	97	THR

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Mol	Chain	Res	Type
27	DK	104	THR
27	DK	105	ARG
27	DK	108	ARG
27	DK	111	LYS
27	DK	113	MET
27	DK	120	PRO
28	DP	3	ILE
28	DP	6	GLN
28	DP	19	PHE
28	DP	20	ARG
28	DP	25	VAL
28	DP	33	GLU
28	DP	43	GLU
28	DP	52	ARG
28	DP	58	PHE
28	DP	59	THR
28	DP	61	ARG
28	DP	65	ASN
28	DP	80	VAL
28	DP	83	ILE
28	DP	84	SER
28	DP	95	LYS
28	DP	99	LEU
28	DP	100	ARG
28	DP	101	GLU
28	DP	111	GLU
28	DP	112	ARG
28	DP	114	ASN
29	DE	5	LEU
29	DE	6	LYS
29	DE	7	ASP
29	DE	9	GLN
29	DE	13	THR
29	DE	16	GLU
29	DE	17	THR
29	DE	24	ASN
29	DE	58	LYS
29	DE	61	ARG
29	DE	62	GLN
29	DE	63	LYS
29	DE	78	TRP
29	DE	108	ILE

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Mol	Chain	Res	Type
29	DE	112	LEU
29	DE	118	LEU
29	DE	122	GLU
29	DE	123	LYS
29	DE	144	GLU
29	DE	145	ASP
29	DE	147	LEU
29	DE	198	GLU
30	DY	2	LYS
30	DY	6	ILE
30	DY	10	ARG
30	DY	15	ARG
30	DY	23	LEU
30	DY	24	LEU
30	DY	37	ARG
30	DY	43	ILE
30	DY	53	MET
31	D0	5	ASN
31	D0	27	LEU
31	D0	28	SER
31	D0	35	GLU
31	D0	37	HIS
31	D0	38	LEU
31	D0	41	HIS
31	D0	45	ASP
31	D0	47	TYR
31	D0	51	ARG
32	D4	9	LYS
32	D4	24	ARG
32	D4	35	GLN
33	D1	6	GLU
33	D1	8	ILE
33	D1	9	LYS
33	D1	16	THR
33	D1	31	GLU
33	D1	35	LEU
34	D3	7	ARG
34	D3	14	LYS
34	D3	18	LYS
34	D3	56	LEU
34	D3	57	VAL
34	D3	58	ILE

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Mol	Chain	Res	Type
35	DV	11	GLU
35	DV	18	ARG
35	DV	24	ASN
35	DV	34	LYS
35	DV	40	ILE
35	DV	42	LEU
35	DV	51	GLN
35	DV	53	LYS
35	DV	68	LYS
35	DV	70	ILE
35	DV	75	GLN
35	DV	77	VAL
35	DV	87	GLN
35	DV	90	ASP
36	D2	3	ARG
36	D2	22	MET
36	D2	34	ARG
36	D2	39	ARG
36	D2	46	LYS
37	DL	4	ASN
37	DL	19	LEU
37	DL	21	ARG
37	DL	46	VAL
37	DL	64	PHE
37	DL	69	ARG
37	DL	92	LEU
37	DL	94	THR
37	DL	95	LEU
37	DL	118	THR
37	DL	122	VAL
37	DL	123	ARG
38	DM	1	MET
38	DM	20	LEU
38	DM	25	ASP
38	DM	42	THR
38	DM	46	ILE
38	DM	50	ARG
38	DM	55	ARG
38	DM	59	ARG
38	DM	60	GLN
38	DM	63	ILE
38	DM	65	ILE

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Mol	Chain	Res	Type
38	DM	70	ASP
38	DM	75	GLU
38	DM	78	LEU
38	DM	81	ARG
38	DM	82	MET
38	DM	88	ASN
38	DM	90	GLU
38	DM	93	VAL
38	DM	110	GLU
38	DM	111	GLU
38	DM	115	GLU
38	DM	131	VAL
39	DX	2	LYS
39	DX	7	ARG
39	DX	16	THR
39	DX	17	GLU
39	DX	20	ASN
39	DX	25	GLN
39	DX	37	LEU
39	DX	57	LEU
39	DX	58	ASN
39	DX	59	GLU
40	DH	1	MET
40	DH	3	VAL
40	DH	12	LEU
40	DH	15	LEU
40	DH	17	ASP
40	DH	25	TYR
40	DH	28	ASN
40	DH	50	ARG
40	DH	70	GLU
40	DH	76	GLU
40	DH	87	GLU
40	DH	90	LEU
40	DH	96	THR
40	DH	97	ARG
40	DH	112	LYS
40	DH	116	ARG
40	DH	119	ASN
40	DH	123	ARG
40	DH	124	THR
40	DH	129	GLU

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Mol	Chain	Res	Type
40	DH	131	SER
40	DH	134	VAL
40	DH	137	GLU
40	DH	139	PHE
40	DH	141	LYS
41	DJ	2	LYS
41	DJ	3	THR
41	DJ	5	THR
41	DJ	12	LYS
41	DJ	21	THR
41	DJ	31	GLU
41	DJ	43	GLU
41	DJ	44	TYR
41	DJ	50	THR
41	DJ	54	ILE
41	DJ	65	THR
41	DJ	72	LYS
41	DJ	73	VAL
41	DJ	95	ARG
41	DJ	106	LYS
41	DJ	108	MET
41	DJ	129	GLU
41	DJ	131	ASN
41	DJ	136	GLN
41	DJ	140	LEU
42	DN	2	ARG
42	DN	3	HIS
42	DN	11	ASN
42	DN	22	ARG
42	DN	33	ILE
42	DN	35	LYS
42	DN	46	ARG
42	DN	51	LEU
42	DN	57	THR
42	DN	69	ARG
42	DN	100	CYS
42	DN	114	GLU
42	DN	120	GLU
43	DO	3	LYS
43	DO	9	ARG
43	DO	21	LEU
43	DO	31	THR

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Mol	Chain	Res	Type
43	DO	33	ARG
43	DO	35	ILE
43	DO	43	ASN
43	DO	46	GLU
43	DO	78	VAL
43	DO	84	GLU
43	DO	98	GLN
43	DO	104	GLN
43	DO	106	LEU
43	DO	116	GLN
44	DQ	5	ARG
44	DQ	10	ARG
44	DQ	13	HIS
44	DQ	14	LYS
44	DQ	19	GLN
44	DQ	29	ARG
44	DQ	48	ASP
44	DQ	59	LEU
44	DQ	69	ARG
44	DQ	79	ILE
44	DQ	83	LYS
44	DQ	88	GLU
44	DQ	90	ASP
44	DQ	96	ASP
44	DQ	105	PHE
45	DS	1	MET
45	DS	6	LYS
45	DS	17	VAL
45	DS	18	ARG
45	DS	22	ASP
45	DS	36	LEU
45	DS	40	ASN
45	DS	66	ILE
45	DS	72	THR
45	DS	84	ARG
45	DS	86	MET
45	DS	88	ARG
45	DS	101	SER
46	DU	7	ASP
46	DU	11	ILE
46	DU	18	LYS
46	DU	26	ASN

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Mol	Chain	Res	Type
46	DU	35	VAL
46	DU	45	GLN
46	DU	46	LYS
46	DU	51	LEU
46	DU	60	LYS
46	DU	73	ASN
46	DU	78	LYS
46	DU	85	ARG
46	DU	87	GLU
46	DU	102	ILE
47	DF	2	LYS
47	DF	9	ASP
47	DF	18	GLU
47	DF	22	ASN
47	DF	29	ARG
47	DF	35	LEU
47	DF	49	LEU
47	DF	50	ASP
47	DF	56	LEU
47	DF	62	GLN
47	DF	66	ILE
47	DF	68	LYS
47	DF	76	PHE
47	DF	91	ARG
47	DF	93	GLU
47	DF	94	ARG
47	DF	97	GLU
47	DF	100	GLU
47	DF	111	ARG
47	DF	121	PHE
47	DF	129	MET
47	DF	134	GLN
47	DF	137	PHE
47	DF	138	PRO
47	DF	141	ASP
47	DF	142	TYR
47	DF	143	ASP
47	DF	146	ASP
47	DF	147	ARG
47	DF	149	ARG
47	DF	156	THR
47	DF	177	ARG

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Mol	Chain	Res	Type
47	DF	178	LYS
48	DG	14	VAL
48	DG	17	LYS
48	DG	19	ASN
48	DG	26	LYS
48	DG	29	ASN
48	DG	31	GLU
48	DG	34	ARG
48	DG	36	LEU
48	DG	46	ASP
48	DG	50	THR
48	DG	54	ARG
48	DG	68	ARG
48	DG	84	LYS
48	DG	85	LYS
48	DG	86	LEU
48	DG	94	ARG
48	DG	120	ILE
48	DG	123	GLU
48	DG	138	GLN
48	DG	148	ARG
48	DG	162	ARG
48	DG	166	GLU
48	DG	167	VAL
48	DG	169	ARG
48	DG	176	LYS
49	DR	5	PHE
49	DR	12	HIS
49	DR	19	THR
49	DR	22	LEU
49	DR	27	ILE
49	DR	39	LEU
49	DR	41	ILE
49	DR	48	LYS
49	DR	53	PHE
49	DR	70	GLU
49	DR	99	THR
50	DT	2	ILE
50	DT	3	ARG
50	DT	6	ARG
50	DT	9	LYS
50	DT	11	LEU

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Mol	Chain	Res	Type
50	DT	12	ARG
50	DT	22	THR
50	DT	29	THR
50	DT	32	LEU
50	DT	50	LEU
50	DT	54	GLU
50	DT	61	LEU
50	DT	64	LYS
50	DT	68	LYS
50	DT	69	ARG
50	DT	73	ARG
50	DT	76	ARG
50	DT	79	ASP
50	DT	87	LEU
50	DT	92	ASN
51	DZ	4	VAL
51	DZ	13	VAL
51	DZ	27	ARG
51	DZ	28	ARG
51	DZ	29	PHE
51	DZ	30	LEU
51	DZ	33	LEU
51	DZ	37	ARG
51	DZ	46	PHE
51	DZ	49	LEU
51	DZ	60	ASP
51	DZ	65	ASP
51	DZ	77	LYS
51	DZ	78	TYR
52	DW	10	ARG
52	DW	16	GLU
52	DW	18	LYS
52	DW	19	ARG
52	DW	23	LYS
52	DW	24	ARG
52	DW	25	PHE
52	DW	31	LEU
52	DW	38	ARG
52	DW	39	GLN
52	DW	40	ARG
52	DW	44	PHE
52	DW	50	VAL

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Mol	Chain	Res	Type
52	DW	61	LYS
52	DW	75	ASN
52	DW	76	ARG
52	DW	77	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (332) such sidechains are listed below:

Mol	Chain	Res	Type
2	AC	2	GLN
2	AC	139	ASN
2	AC	184	ASN
3	AD	35	GLN
3	AD	53	GLN
3	AD	115	GLN
3	AD	119	HIS
3	AD	135	GLN
3	AD	151	GLN
3	AD	195	ASN
4	AE	18	ASN
4	AE	42	ASN
4	AE	72	ASN
4	AE	77	ASN
4	AE	81	GLN
4	AE	82	HIS
4	AE	131	ASN
4	AE	134	ASN
5	AF	17	GLN
5	AF	46	GLN
5	AF	58	HIS
6	AG	67	ASN
6	AG	85	GLN
6	AG	121	ASN
6	AG	129	ASN
6	AG	141	HIS
6	AG	147	ASN
7	AH	3	GLN
7	AH	17	GLN
7	AH	75	GLN
7	AH	117	GLN
8	AI	30	ASN
8	AI	31	GLN
8	AI	36	GLN

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Mol	Chain	Res	Type
8	AI	109	GLN
9	AJ	35	GLN
9	AJ	58	ASN
10	AK	21	HIS
10	AK	28	ASN
10	AK	39	ASN
10	AK	80	ASN
10	AK	108	ASN
10	AK	118	ASN
11	AL	19	ASN
11	AL	28	GLN
11	AL	45	ASN
11	AL	111	GLN
12	AM	90	HIS
13	AP	29	ASN
13	AP	40	ASN
13	AP	59	HIS
14	AQ	8	GLN
17	AT	47	GLN
18	AB	23	ASN
18	AB	35	ASN
18	AB	121	GLN
18	AB	145	ASN
18	AB	169	HIS
18	AB	177	ASN
18	AB	202	ASN
19	AU	8	ASN
20	AO	20	ASN
20	AO	37	ASN
20	AO	40	GLN
21	AN	59	GLN
21	AN	65	GLN
21	AN	70	HIS
24	BI	11	GLN
24	BI	29	GLN
24	BI	33	ASN
24	BI	93	ASN
25	BC	20	ASN
25	BC	36	ASN
25	BC	43	ASN
25	BC	52	HIS
25	BC	59	GLN

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Mol	Chain	Res	Type
25	BC	89	ASN
25	BC	152	GLN
25	BC	162	GLN
25	BC	225	ASN
26	BD	32	ASN
26	BD	36	GLN
26	BD	49	GLN
26	BD	126	ASN
26	BD	136	ASN
27	BK	5	GLN
27	BK	88	ASN
27	BK	89	ASN
27	BK	93	GLN
28	BP	6	GLN
28	BP	114	ASN
29	BE	9	GLN
29	BE	24	ASN
29	BE	29	HIS
29	BE	30	GLN
29	BE	62	GLN
29	BE	156	ASN
29	BE	165	HIS
29	BE	195	GLN
30	BY	48	ASN
32	B4	13	ASN
32	B4	37	GLN
33	B1	25	ASN
34	B3	30	HIS
35	BV	44	HIS
35	BV	49	ASN
35	BV	51	GLN
35	BV	80	HIS
35	BV	88	HIS
36	B2	6	GLN
36	B2	13	ASN
36	B2	26	ASN
36	B2	29	GLN
37	BL	4	ASN
37	BL	54	GLN
37	BL	93	ASN
37	BL	104	GLN
38	BM	3	GLN

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Mol	Chain	Res	Type
38	BM	17	ASN
38	BM	45	GLN
39	BX	15	ASN
39	BX	20	ASN
39	BX	25	GLN
39	BX	27	ASN
39	BX	31	GLN
39	BX	58	ASN
40	BH	18	GLN
40	BH	28	ASN
40	BH	66	ASN
40	BH	73	ASN
41	BJ	130	HIS
42	BN	11	ASN
42	BN	23	ASN
42	BN	62	ASN
42	BN	107	ASN
43	BO	19	GLN
43	BO	38	GLN
43	BO	61	GLN
43	BO	116	GLN
44	BQ	55	GLN
44	BQ	58	GLN
44	BQ	80	ASN
45	BS	57	ASN
46	BU	26	ASN
46	BU	53	GLN
46	BU	68	ASN
46	BU	73	ASN
47	BF	22	ASN
47	BF	51	ASN
47	BF	62	GLN
47	BF	126	ASN
47	BF	134	GLN
48	BG	29	ASN
48	BG	63	GLN
48	BG	127	GLN
49	BR	6	GLN
49	BR	12	HIS
49	BR	86	GLN
50	BT	48	GLN
50	BT	92	ASN

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Mol	Chain	Res	Type
51	BZ	6	GLN
51	BZ	20	HIS
51	BZ	23	ASN
51	BZ	32	ASN
51	BZ	36	HIS
52	BW	11	ASN
52	BW	56	HIS
52	BW	75	ASN
3	CD	35	GLN
3	CD	53	GLN
3	CD	73	ASN
3	CD	84	ASN
3	CD	115	GLN
3	CD	119	HIS
3	CD	135	GLN
3	CD	151	GLN
3	CD	195	ASN
4	CE	18	ASN
4	CE	42	ASN
4	CE	72	ASN
4	CE	77	ASN
4	CE	81	GLN
4	CE	82	HIS
4	CE	131	ASN
4	CE	134	ASN
5	CF	17	GLN
5	CF	46	GLN
5	CF	58	HIS
7	CH	3	GLN
7	CH	17	GLN
7	CH	75	GLN
7	CH	117	GLN
10	CK	28	ASN
10	CK	39	ASN
10	CK	80	ASN
10	CK	108	ASN
10	CK	118	ASN
11	CL	19	ASN
11	CL	28	GLN
11	CL	45	ASN
11	CL	111	GLN
13	CP	29	ASN

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Mol	Chain	Res	Type
13	CP	40	ASN
13	CP	59	HIS
14	CQ	8	GLN
17	CT	47	GLN
2	CC	2	GLN
2	CC	5	HIS
2	CC	31	ASN
2	CC	68	HIS
2	CC	101	ASN
2	CC	139	ASN
2	CC	184	ASN
2	CC	189	HIS
6	CG	67	ASN
6	CG	152	HIS
8	CI	30	ASN
8	CI	31	GLN
8	CI	36	GLN
8	CI	125	GLN
9	CJ	20	GLN
9	CJ	64	GLN
9	CJ	99	GLN
12	CM	13	HIS
21	CN	59	GLN
21	CN	65	GLN
16	CS	42	ASN
18	CB	14	HIS
18	CB	35	ASN
18	CB	88	GLN
18	CB	108	GLN
18	CB	119	GLN
18	CB	145	ASN
18	CB	169	HIS
20	CO	35	GLN
20	CO	38	HIS
24	DI	5	GLN
24	DI	11	GLN
24	DI	33	ASN
25	DC	20	ASN
25	DC	36	ASN
25	DC	43	ASN
25	DC	59	GLN
25	DC	89	ASN

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Mol	Chain	Res	Type
25	DC	152	GLN
25	DC	162	GLN
25	DC	196	ASN
26	DD	32	ASN
26	DD	36	GLN
26	DD	49	GLN
26	DD	126	ASN
26	DD	136	ASN
27	DK	5	GLN
27	DK	88	ASN
27	DK	89	ASN
27	DK	93	GLN
28	DP	6	GLN
28	DP	74	GLN
28	DP	114	ASN
29	DE	9	GLN
29	DE	24	ASN
29	DE	29	HIS
29	DE	30	GLN
29	DE	62	GLN
29	DE	156	ASN
29	DE	165	HIS
29	DE	195	GLN
30	DY	48	ASN
32	D4	13	ASN
32	D4	37	GLN
34	D3	30	HIS
35	DV	44	HIS
35	DV	51	GLN
35	DV	80	HIS
36	D2	6	GLN
36	D2	13	ASN
36	D2	29	GLN
37	DL	4	ASN
37	DL	54	GLN
37	DL	93	ASN
37	DL	104	GLN
38	DM	17	ASN
38	DM	45	GLN
39	DX	15	ASN
39	DX	20	ASN
39	DX	25	GLN

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Mol	Chain	Res	Type
39	DX	27	ASN
39	DX	31	GLN
39	DX	58	ASN
40	DH	18	GLN
40	DH	28	ASN
40	DH	66	ASN
40	DH	119	ASN
41	DJ	130	HIS
42	DN	11	ASN
42	DN	23	ASN
42	DN	62	ASN
42	DN	107	ASN
43	DO	19	GLN
43	DO	38	GLN
43	DO	61	GLN
43	DO	116	GLN
44	DQ	55	GLN
44	DQ	58	GLN
44	DQ	80	ASN
45	DS	57	ASN
46	DU	26	ASN
46	DU	53	GLN
46	DU	65	GLN
46	DU	68	ASN
46	DU	73	ASN
47	DF	22	ASN
47	DF	51	ASN
47	DF	62	GLN
47	DF	126	ASN
47	DF	134	GLN
48	DG	29	ASN
48	DG	127	GLN
48	DG	142	GLN
49	DR	6	GLN
49	DR	12	HIS
49	DR	86	GLN
50	DT	48	GLN
50	DT	92	ASN
51	DZ	6	GLN
51	DZ	20	HIS
51	DZ	23	ASN
51	DZ	32	ASN

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Mol	Chain	Res	Type
51	DZ	36	HIS
52	DW	11	ASN
52	DW	56	HIS
52	DW	75	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1529/1542 (99%)	292 (19%)	25 (1%)
1	CA	1529/1542 (99%)	282 (18%)	21 (1%)
22	BA	116/120 (96%)	21 (18%)	1 (0%)
22	DA	116/120 (96%)	19 (16%)	1 (0%)
23	BB	2837/2904 (97%)	456 (16%)	18 (0%)
23	DB	2837/2904 (97%)	469 (16%)	20 (0%)
All	All	8964/9132 (98%)	1539 (17%)	86 (0%)

All (1539) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	7	A
1	AA	9	G
1	AA	14	U
1	AA	32	A
1	AA	39	G
1	AA	47	C
1	AA	48	C
1	AA	49	U
1	AA	52	C
1	AA	55	A
1	AA	61	G
1	AA	65	A
1	AA	66	A
1	AA	72	A
1	AA	75	G
1	AA	78	A
1	AA	79	G
1	AA	80	A
1	AA	81	A
1	AA	82	G
1	AA	83	C
1	AA	84	U

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Mol	Chain	Res	Type
1	AA	85	U
1	AA	86	G
1	AA	88	U
1	AA	89	U
1	AA	91	U
1	AA	92	U
1	AA	94	G
1	AA	95	C
1	AA	108	G
1	AA	120	A
1	AA	121	U
1	AA	122	G
1	AA	131	A
1	AA	149	A
1	AA	151	A
1	AA	164	G
1	AA	177	G
1	AA	180	U
1	AA	182	A
1	AA	196	A
1	AA	197	A
1	AA	209	U
1	AA	210	C
1	AA	213	G
1	AA	226	G
1	AA	233	C
1	AA	239	U
1	AA	240	G
1	AA	243	A
1	AA	244	U
1	AA	245	U
1	AA	247	G
1	AA	250	A
1	AA	251	G
1	AA	253	A
1	AA	256	U
1	AA	257	G
1	AA	258	G
1	AA	266	G
1	AA	267	C
1	AA	280	C
1	AA	289	G

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Mol	Chain	Res	Type
1	AA	301	G
1	AA	316	C
1	AA	321	A
1	AA	328	C
1	AA	329	A
1	AA	330	C
1	AA	332	G
1	AA	345	C
1	AA	352	C
1	AA	354	G
1	AA	367	U
1	AA	372	C
1	AA	373	A
1	AA	381	C
1	AA	382	A
1	AA	397	A
1	AA	398	U
1	AA	406	G
1	AA	408	A
1	AA	409	U
1	AA	410	G
1	AA	411	A
1	AA	412	A
1	AA	413	G
1	AA	414	A
1	AA	415	A
1	AA	421	U
1	AA	422	C
1	AA	429	U
1	AA	430	A
1	AA	435	A
1	AA	438	U
1	AA	457	G
1	AA	461	A
1	AA	462	G
1	AA	463	U
1	AA	464	U
1	AA	465	A
1	AA	466	A
1	AA	467	U
1	AA	468	A
1	AA	479	U

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Mol	Chain	Res	Type
1	AA	484	G
1	AA	485	U
1	AA	486	U
1	AA	493	A
1	AA	500	G
1	AA	511	C
1	AA	518	C
1	AA	521	G
1	AA	524	G
1	AA	527	G
1	AA	532	A
1	AA	547	A
1	AA	559	A
1	AA	562	U
1	AA	566	G
1	AA	572	A
1	AA	573	A
1	AA	576	C
1	AA	577	G
1	AA	616	G
1	AA	633	G
1	AA	639	G
1	AA	653	U
1	AA	665	A
1	AA	688	G
1	AA	700	G
1	AA	703	G
1	AA	721	G
1	AA	724	G
1	AA	731	G
1	AA	733	G
1	AA	747	A
1	AA	748	G
1	AA	755	G
1	AA	777	A
1	AA	781	A
1	AA	782	A
1	AA	793	U
1	AA	794	A
1	AA	812	G
1	AA	813	U
1	AA	815	A

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Mol	Chain	Res	Type
1	AA	817	C
1	AA	819	A
1	AA	828	U
1	AA	829	G
1	AA	841	C
1	AA	842	U
1	AA	843	U
1	AA	844	G
1	AA	845	A
1	AA	847	G
1	AA	858	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	968	A
1	AA	969	A
1	AA	971	G
1	AA	974	A
1	AA	976	G
1	AA	977	A
1	AA	991	U
1	AA	992	U
1	AA	993	G
1	AA	994	A
1	AA	995	C
1	AA	1000	A
1	AA	1002	G
1	AA	1003	G
1	AA	1004	A
1	AA	1010	U
1	AA	1015	G
1	AA	1020	G
1	AA	1022	A
1	AA	1026	G
1	AA	1028	C
1	AA	1030	U
1	AA	1031	C
1	AA	1033	G
1	AA	1041	G

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Mol	Chain	Res	Type
1	AA	1044	A
1	AA	1045	C
1	AA	1049	U
1	AA	1050	G
1	AA	1051	C
1	AA	1054	C
1	AA	1055	A
1	AA	1056	U
1	AA	1063	C
1	AA	1064	G
1	AA	1065	U
1	AA	1066	C
1	AA	1094	G
1	AA	1095	U
1	AA	1101	A
1	AA	1102	A
1	AA	1111	A
1	AA	1112	C
1	AA	1118	U
1	AA	1126	U
1	AA	1136	C
1	AA	1137	C
1	AA	1138	G
1	AA	1139	G
1	AA	1140	C
1	AA	1159	U
1	AA	1167	A
1	AA	1168	U
1	AA	1171	A
1	AA	1174	G
1	AA	1181	G
1	AA	1182	G
1	AA	1183	U
1	AA	1196	A
1	AA	1197	A
1	AA	1202	U
1	AA	1209	C
1	AA	1212	U
1	AA	1213	A
1	AA	1214	C
1	AA	1215	G
1	AA	1225	A

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Mol	Chain	Res	Type
1	AA	1226	C
1	AA	1227	A
1	AA	1228	C
1	AA	1233	G
1	AA	1239	A
1	AA	1240	U
1	AA	1241	G
1	AA	1256	A
1	AA	1278	G
1	AA	1279	G
1	AA	1280	A
1	AA	1281	C
1	AA	1286	U
1	AA	1287	A
1	AA	1294	G
1	AA	1297	G
1	AA	1298	U
1	AA	1300	G
1	AA	1301	U
1	AA	1305	G
1	AA	1316	G
1	AA	1317	C
1	AA	1318	A
1	AA	1319	A
1	AA	1320	C
1	AA	1323	G
1	AA	1324	A
1	AA	1334	G
1	AA	1336	C
1	AA	1345	U
1	AA	1347	G
1	AA	1353	G
1	AA	1362	A
1	AA	1364	U
1	AA	1378	C
1	AA	1379	G
1	AA	1380	U
1	AA	1381	U
1	AA	1397	C
1	AA	1398	A
1	AA	1407	C
1	AA	1409	C

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Mol	Chain	Res	Type
1	AA	1411	C
1	AA	1432	G
1	AA	1446	A
1	AA	1448	C
1	AA	1452	C
1	AA	1493	A
1	AA	1497	G
1	AA	1499	A
1	AA	1503	A
1	AA	1504	G
1	AA	1506	U
1	AA	1517	G
1	AA	1519	A
1	AA	1520	C
1	AA	1529	G
1	AA	1530	G
1	AA	1533	C
1	AA	1534	A
22	BA	4	C
22	BA	12	C
22	BA	13	G
22	BA	14	U
22	BA	15	A
22	BA	16	G
22	BA	24	G
22	BA	25	U
22	BA	26	C
22	BA	29	A
22	BA	30	C
22	BA	44	G
22	BA	52	A
22	BA	55	U
22	BA	67	G
22	BA	74	U
22	BA	88	C
22	BA	89	U
22	BA	90	C
22	BA	99	A
22	BA	109	A
23	BB	4	U
23	BB	27	G
23	BB	34	U

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Mol	Chain	Res	Type
23	BB	46	G
23	BB	49	A
23	BB	71	A
23	BB	72	U
23	BB	74	A
23	BB	75	G
23	BB	79	C
23	BB	96	C
23	BB	100	U
23	BB	101	A
23	BB	102	U
23	BB	109	C
23	BB	110	G
23	BB	118	A
23	BB	119	A
23	BB	120	U
23	BB	121	G
23	BB	125	A
23	BB	136	G
23	BB	137	U
23	BB	138	U
23	BB	139	U
23	BB	140	C
23	BB	141	G
23	BB	142	A
23	BB	144	A
23	BB	160	A
23	BB	163	C
23	BB	180	G
23	BB	181	A
23	BB	196	A
23	BB	199	A
23	BB	215	G
23	BB	216	A
23	BB	221	A
23	BB	222	A
23	BB	223	A
23	BB	233	A
23	BB	241	A
23	BB	248	G
23	BB	249	C
23	BB	255	A

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Mol	Chain	Res	Type
23	BB	264	C
23	BB	265	A
23	BB	266	G
23	BB	268	C
23	BB	271	G
23	BB	277	G
23	BB	279	A
23	BB	281	C
23	BB	294	A
23	BB	299	A
23	BB	311	A
23	BB	322	A
23	BB	323	C
23	BB	329	G
23	BB	330	A
23	BB	331	C
23	BB	333	G
23	BB	341	C
23	BB	343	C
23	BB	346	A
23	BB	349	U
23	BB	353	C
23	BB	355	U
23	BB	371	A
23	BB	372	G
23	BB	386	G
23	BB	387	U
23	BB	396	G
23	BB	405	U
23	BB	406	G
23	BB	411	G
23	BB	424	G
23	BB	451	U
23	BB	455	C
23	BB	457	A
23	BB	479	A
23	BB	481	G
23	BB	483	A
23	BB	490	C
23	BB	491	G
23	BB	504	A
23	BB	505	A

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Mol	Chain	Res	Type
23	BB	508	A
23	BB	509	C
23	BB	512	G
23	BB	527	C
23	BB	532	A
23	BB	533	G
23	BB	542	C
23	BB	544	C
23	BB	545	U
23	BB	546	U
23	BB	547	A
23	BB	548	G
23	BB	549	G
23	BB	550	C
23	BB	563	A
23	BB	568	U
23	BB	573	U
23	BB	574	A
23	BB	575	A
23	BB	603	A
23	BB	613	A
23	BB	614	A
23	BB	615	U
23	BB	627	A
23	BB	628	G
23	BB	637	A
23	BB	645	C
23	BB	646	U
23	BB	653	U
23	BB	654	A
23	BB	655	A
23	BB	656	G
23	BB	671	C
23	BB	676	A
23	BB	686	U
23	BB	699	A
23	BB	730	A
23	BB	747	U
23	BB	757	G
23	BB	775	G
23	BB	776	G
23	BB	782	A

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Mol	Chain	Res	Type
23	BB	784	G
23	BB	785	G
23	BB	802	A
23	BB	805	G
23	BB	806	C
23	BB	812	C
23	BB	819	A
23	BB	827	U
23	BB	828	U
23	BB	859	G
23	BB	871	U
23	BB	872	U
23	BB	877	A
23	BB	899	A
23	BB	910	A
23	BB	912	C
23	BB	931	U
23	BB	932	U
23	BB	933	A
23	BB	941	A
23	BB	946	C
23	BB	958	U
23	BB	961	C
23	BB	973	A
23	BB	974	G
23	BB	980	A
23	BB	982	C
23	BB	983	A
23	BB	990	A
23	BB	991	C
23	BB	995	C
23	BB	996	A
23	BB	1005	C
23	BB	1012	U
23	BB	1013	C
23	BB	1022	G
23	BB	1023	U
23	BB	1024	G
23	BB	1025	G
23	BB	1033	U
23	BB	1042	G
23	BB	1044	C

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Mol	Chain	Res	Type
23	BB	1047	G
23	BB	1062	G
23	BB	1070	A
23	BB	1088	A
23	BB	1090	A
23	BB	1110	G
23	BB	1112	G
23	BB	1115	G
23	BB	1129	A
23	BB	1132	U
23	BB	1133	A
23	BB	1134	A
23	BB	1135	C
23	BB	1136	G
23	BB	1141	U
23	BB	1157	G
23	BB	1172	C
23	BB	1173	U
23	BB	1174	U
23	BB	1175	A
23	BB	1176	U
23	BB	1195	G
23	BB	1204	A
23	BB	1206	G
23	BB	1211	C
23	BB	1212	G
23	BB	1237	A
23	BB	1241	A
23	BB	1242	U
23	BB	1248	G
23	BB	1250	G
23	BB	1251	C
23	BB	1253	A
23	BB	1256	G
23	BB	1266	G
23	BB	1271	G
23	BB	1272	A
23	BB	1273	U
23	BB	1275	A
23	BB	1300	G
23	BB	1301	A
23	BB	1318	U

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Mol	Chain	Res	Type
23	BB	1321	A
23	BB	1324	G
23	BB	1325	U
23	BB	1337	G
23	BB	1341	G
23	BB	1365	A
23	BB	1368	G
23	BB	1379	U
23	BB	1383	A
23	BB	1384	A
23	BB	1386	C
23	BB	1396	U
23	BB	1397	U
23	BB	1416	G
23	BB	1419	A
23	BB	1421	G
23	BB	1427	A
23	BB	1428	C
23	BB	1434	A
23	BB	1453	A
23	BB	1454	C
23	BB	1458	U
23	BB	1459	G
23	BB	1460	U
23	BB	1461	C
23	BB	1476	U
23	BB	1477	A
23	BB	1478	G
23	BB	1482	G
23	BB	1490	A
23	BB	1493	C
23	BB	1494	A
23	BB	1504	A
23	BB	1505	A
23	BB	1506	U
23	BB	1507	C
23	BB	1508	A
23	BB	1509	A
23	BB	1523	U
23	BB	1524	G
23	BB	1530	G
23	BB	1532	A

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Mol	Chain	Res	Type
23	BB	1535	A
23	BB	1536	C
23	BB	1552	A
23	BB	1559	U
23	BB	1560	G
23	BB	1567	G
23	BB	1569	A
23	BB	1578	U
23	BB	1583	A
23	BB	1585	C
23	BB	1602	U
23	BB	1603	A
23	BB	1607	C
23	BB	1608	A
23	BB	1610	A
23	BB	1613	G
23	BB	1616	A
23	BB	1618	A
23	BB	1619	G
23	BB	1634	A
23	BB	1635	A
23	BB	1640	A
23	BB	1647	U
23	BB	1648	U
23	BB	1670	C
23	BB	1674	G
23	BB	1681	G
23	BB	1700	A
23	BB	1701	A
23	BB	1715	G
23	BB	1722	A
23	BB	1724	G
23	BB	1727	C
23	BB	1729	U
23	BB	1730	C
23	BB	1731	G
23	BB	1733	G
23	BB	1738	G
23	BB	1756	G
23	BB	1758	U
23	BB	1764	C
23	BB	1773	A

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Mol	Chain	Res	Type
23	BB	1800	C
23	BB	1801	A
23	BB	1808	A
23	BB	1816	C
23	BB	1848	A
23	BB	1869	G
23	BB	1870	C
23	BB	1876	A
23	BB	1884	G
23	BB	1906	G
23	BB	1914	C
23	BB	1915	U
23	BB	1926	U
23	BB	1929	G
23	BB	1930	G
23	BB	1937	A
23	BB	1938	A
23	BB	1939	U
23	BB	1940	U
23	BB	1955	U
23	BB	1964	G
23	BB	1966	A
23	BB	1967	C
23	BB	1970	A
23	BB	1971	U
23	BB	1972	G
23	BB	1991	U
23	BB	1992	G
23	BB	1993	U
23	BB	1996	C
23	BB	1997	C
23	BB	2022	U
23	BB	2023	C
23	BB	2030	A
23	BB	2031	A
23	BB	2032	G
23	BB	2043	C
23	BB	2055	C
23	BB	2056	G
23	BB	2060	A
23	BB	2061	G
23	BB	2062	A

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Mol	Chain	Res	Type
23	BB	2069	G
23	BB	2072	C
23	BB	2093	G
23	BB	2095	A
23	BB	2134	A
23	BB	2141	G
23	BB	2142	A
23	BB	2144	G
23	BB	2145	C
23	BB	2147	A
23	BB	2150	C
23	BB	2154	A
23	BB	2180	U
23	BB	2181	U
23	BB	2182	U
23	BB	2184	A
23	BB	2192	U
23	BB	2198	A
23	BB	2199	A
23	BB	2203	U
23	BB	2204	G
23	BB	2210	U
23	BB	2211	A
23	BB	2212	A
23	BB	2213	U
23	BB	2214	C
23	BB	2225	A
23	BB	2238	G
23	BB	2239	G
23	BB	2249	U
23	BB	2250	G
23	BB	2279	G
23	BB	2283	C
23	BB	2287	A
23	BB	2288	A
23	BB	2304	G
23	BB	2305	U
23	BB	2307	G
23	BB	2308	G
23	BB	2309	A
23	BB	2310	C
23	BB	2311	A

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Mol	Chain	Res	Type
23	BB	2321	U
23	BB	2322	A
23	BB	2324	U
23	BB	2325	G
23	BB	2333	A
23	BB	2336	A
23	BB	2337	G
23	BB	2345	G
23	BB	2347	C
23	BB	2357	G
23	BB	2361	G
23	BB	2372	U
23	BB	2383	G
23	BB	2385	C
23	BB	2396	G
23	BB	2402	U
23	BB	2403	C
23	BB	2406	A
23	BB	2423	U
23	BB	2426	A
23	BB	2429	G
23	BB	2430	A
23	BB	2431	U
23	BB	2441	U
23	BB	2448	A
23	BB	2472	G
23	BB	2476	A
23	BB	2491	U
23	BB	2492	U
23	BB	2502	G
23	BB	2505	G
23	BB	2506	U
23	BB	2518	A
23	BB	2529	G
23	BB	2543	G
23	BB	2552	U
23	BB	2554	U
23	BB	2566	A
23	BB	2567	G
23	BB	2572	A
23	BB	2573	C
23	BB	2574	G

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Mol	Chain	Res	Type
23	BB	2586	U
23	BB	2602	A
23	BB	2609	U
23	BB	2613	U
23	BB	2629	U
23	BB	2639	A
23	BB	2682	A
23	BB	2689	U
23	BB	2690	U
23	BB	2714	G
23	BB	2737	G
23	BB	2744	G
23	BB	2757	A
23	BB	2765	A
23	BB	2778	A
23	BB	2791	G
23	BB	2796	U
23	BB	2798	U
23	BB	2799	A
23	BB	2808	G
23	BB	2820	A
23	BB	2821	A
23	BB	2823	A
23	BB	2835	A
23	BB	2850	A
23	BB	2859	G
23	BB	2867	G
23	BB	2872	A
23	BB	2873	A
23	BB	2880	C
23	BB	2883	A
23	BB	2894	G
23	BB	2903	U
1	CA	7	A
1	CA	9	G
1	CA	14	U
1	CA	32	A
1	CA	39	G
1	CA	47	C
1	CA	48	C
1	CA	49	U
1	CA	52	C

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Mol	Chain	Res	Type
1	CA	54	C
1	CA	55	A
1	CA	61	G
1	CA	71	A
1	CA	72	A
1	CA	74	A
1	CA	76	G
1	CA	83	C
1	CA	85	U
1	CA	86	G
1	CA	87	C
1	CA	88	U
1	CA	92	U
1	CA	93	U
1	CA	94	G
1	CA	108	G
1	CA	120	A
1	CA	121	U
1	CA	122	G
1	CA	131	A
1	CA	149	A
1	CA	151	A
1	CA	164	G
1	CA	177	G
1	CA	180	U
1	CA	182	A
1	CA	196	A
1	CA	197	A
1	CA	209	U
1	CA	210	C
1	CA	213	G
1	CA	226	G
1	CA	233	C
1	CA	239	U
1	CA	240	G
1	CA	243	A
1	CA	244	U
1	CA	245	U
1	CA	247	G
1	CA	250	A
1	CA	251	G
1	CA	253	A

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Mol	Chain	Res	Type
1	CA	254	G
1	CA	256	U
1	CA	257	G
1	CA	258	G
1	CA	266	G
1	CA	267	C
1	CA	280	C
1	CA	289	G
1	CA	301	G
1	CA	316	C
1	CA	321	A
1	CA	328	C
1	CA	329	A
1	CA	330	C
1	CA	332	G
1	CA	345	C
1	CA	352	C
1	CA	354	G
1	CA	367	U
1	CA	372	C
1	CA	373	A
1	CA	381	C
1	CA	382	A
1	CA	397	A
1	CA	398	U
1	CA	406	G
1	CA	408	A
1	CA	409	U
1	CA	410	G
1	CA	411	A
1	CA	412	A
1	CA	413	G
1	CA	414	A
1	CA	415	A
1	CA	421	U
1	CA	422	C
1	CA	429	U
1	CA	430	A
1	CA	435	A
1	CA	438	U
1	CA	457	G
1	CA	461	A

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Mol	Chain	Res	Type
1	CA	462	G
1	CA	463	U
1	CA	464	U
1	CA	465	A
1	CA	466	A
1	CA	467	U
1	CA	468	A
1	CA	479	U
1	CA	482	A
1	CA	484	G
1	CA	485	U
1	CA	486	U
1	CA	493	A
1	CA	500	G
1	CA	511	C
1	CA	518	C
1	CA	521	G
1	CA	524	G
1	CA	527	G
1	CA	532	A
1	CA	547	A
1	CA	559	A
1	CA	562	U
1	CA	566	G
1	CA	572	A
1	CA	573	A
1	CA	576	C
1	CA	577	G
1	CA	616	G
1	CA	633	G
1	CA	639	G
1	CA	653	U
1	CA	665	A
1	CA	688	G
1	CA	700	G
1	CA	703	G
1	CA	721	G
1	CA	724	G
1	CA	731	G
1	CA	733	G
1	CA	747	A
1	CA	748	G

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Mol	Chain	Res	Type
1	CA	755	G
1	CA	777	A
1	CA	781	A
1	CA	782	A
1	CA	793	U
1	CA	794	A
1	CA	812	G
1	CA	813	U
1	CA	815	A
1	CA	817	C
1	CA	819	A
1	CA	828	U
1	CA	829	G
1	CA	841	C
1	CA	842	U
1	CA	843	U
1	CA	844	G
1	CA	845	A
1	CA	847	G
1	CA	858	G
1	CA	914	A
1	CA	926	G
1	CA	934	C
1	CA	935	A
1	CA	958	A
1	CA	960	U
1	CA	961	U
1	CA	966	G
1	CA	968	A
1	CA	969	A
1	CA	971	G
1	CA	972	C
1	CA	974	A
1	CA	975	A
1	CA	976	G
1	CA	977	A
1	CA	991	U
1	CA	992	U
1	CA	993	G
1	CA	1000	A
1	CA	1001	C
1	CA	1004	A

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Mol	Chain	Res	Type
1	CA	1007	U
1	CA	1009	U
1	CA	1019	A
1	CA	1024	G
1	CA	1025	U
1	CA	1026	G
1	CA	1030	U
1	CA	1031	C
1	CA	1032	G
1	CA	1033	G
1	CA	1036	A
1	CA	1041	G
1	CA	1050	G
1	CA	1053	G
1	CA	1054	C
1	CA	1055	A
1	CA	1056	U
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	1125	U
1	CA	1127	G
1	CA	1128	C
1	CA	1135	U
1	CA	1136	C
1	CA	1137	C
1	CA	1138	G
1	CA	1139	G
1	CA	1140	C
1	CA	1145	A
1	CA	1158	C
1	CA	1159	U
1	CA	1160	G
1	CA	1167	A
1	CA	1168	U
1	CA	1174	G
1	CA	1183	U

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Mol	Chain	Res	Type
1	CA	1184	G
1	CA	1196	A
1	CA	1198	G
1	CA	1201	A
1	CA	1202	U
1	CA	1212	U
1	CA	1213	A
1	CA	1214	C
1	CA	1215	G
1	CA	1225	A
1	CA	1226	C
1	CA	1227	A
1	CA	1230	C
1	CA	1240	U
1	CA	1241	G
1	CA	1252	A
1	CA	1256	A
1	CA	1257	A
1	CA	1258	G
1	CA	1280	A
1	CA	1281	C
1	CA	1285	A
1	CA	1286	U
1	CA	1287	A
1	CA	1289	A
1	CA	1290	G
1	CA	1299	A
1	CA	1300	G
1	CA	1301	U
1	CA	1302	C
1	CA	1305	G
1	CA	1316	G
1	CA	1317	C
1	CA	1320	C
1	CA	1323	G
1	CA	1346	A
1	CA	1353	G
1	CA	1357	A
1	CA	1359	C
1	CA	1362	A
1	CA	1363	A
1	CA	1365	G

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Mol	Chain	Res	Type
1	CA	1379	G
1	CA	1381	U
1	CA	1397	C
1	CA	1398	A
1	CA	1419	G
1	CA	1424	U
1	CA	1432	G
1	CA	1446	A
1	CA	1448	C
1	CA	1452	C
1	CA	1493	A
1	CA	1497	G
1	CA	1499	A
1	CA	1503	A
1	CA	1506	U
1	CA	1517	G
1	CA	1519	A
1	CA	1520	C
1	CA	1529	G
1	CA	1530	G
1	CA	1534	A
22	DA	4	C
22	DA	12	C
22	DA	13	G
22	DA	14	U
22	DA	15	A
22	DA	16	G
22	DA	24	G
22	DA	25	U
22	DA	26	C
22	DA	29	A
22	DA	30	C
22	DA	44	G
22	DA	52	A
22	DA	55	U
22	DA	67	G
22	DA	88	C
22	DA	89	U
22	DA	99	A
22	DA	109	A
23	DB	10	A
23	DB	27	G

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Mol	Chain	Res	Type
23	DB	34	U
23	DB	46	G
23	DB	49	A
23	DB	71	A
23	DB	72	U
23	DB	74	A
23	DB	75	G
23	DB	79	C
23	DB	96	C
23	DB	100	U
23	DB	101	A
23	DB	102	U
23	DB	103	A
23	DB	110	G
23	DB	118	A
23	DB	119	A
23	DB	120	U
23	DB	125	A
23	DB	126	A
23	DB	127	A
23	DB	139	U
23	DB	141	G
23	DB	142	A
23	DB	144	A
23	DB	160	A
23	DB	163	C
23	DB	180	G
23	DB	181	A
23	DB	196	A
23	DB	199	A
23	DB	215	G
23	DB	216	A
23	DB	221	A
23	DB	222	A
23	DB	223	A
23	DB	233	A
23	DB	241	A
23	DB	242	G
23	DB	248	G
23	DB	249	C
23	DB	255	A
23	DB	264	C

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Mol	Chain	Res	Type
23	DB	265	A
23	DB	266	G
23	DB	268	C
23	DB	271	G
23	DB	277	G
23	DB	278	A
23	DB	279	A
23	DB	281	C
23	DB	283	G
23	DB	284	U
23	DB	286	U
23	DB	288	U
23	DB	294	A
23	DB	299	A
23	DB	311	A
23	DB	322	A
23	DB	323	C
23	DB	329	G
23	DB	330	A
23	DB	331	C
23	DB	333	G
23	DB	341	C
23	DB	343	C
23	DB	346	A
23	DB	349	U
23	DB	353	C
23	DB	367	G
23	DB	371	A
23	DB	372	G
23	DB	386	G
23	DB	387	U
23	DB	396	G
23	DB	405	U
23	DB	406	G
23	DB	411	G
23	DB	424	G
23	DB	455	C
23	DB	457	A
23	DB	479	A
23	DB	481	G
23	DB	483	A
23	DB	490	C

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Mol	Chain	Res	Type
23	DB	491	G
23	DB	504	A
23	DB	505	A
23	DB	508	A
23	DB	509	C
23	DB	512	G
23	DB	527	C
23	DB	531	C
23	DB	532	A
23	DB	533	G
23	DB	542	C
23	DB	545	U
23	DB	546	U
23	DB	547	A
23	DB	548	G
23	DB	549	G
23	DB	550	C
23	DB	563	A
23	DB	568	U
23	DB	573	U
23	DB	575	A
23	DB	603	A
23	DB	613	A
23	DB	614	A
23	DB	615	U
23	DB	627	A
23	DB	628	G
23	DB	637	A
23	DB	645	C
23	DB	646	U
23	DB	653	U
23	DB	654	A
23	DB	655	A
23	DB	656	G
23	DB	671	C
23	DB	676	A
23	DB	686	U
23	DB	699	A
23	DB	730	A
23	DB	747	U
23	DB	757	G
23	DB	775	G

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Mol	Chain	Res	Type
23	DB	776	G
23	DB	782	A
23	DB	784	G
23	DB	785	G
23	DB	802	A
23	DB	805	G
23	DB	806	C
23	DB	812	C
23	DB	819	A
23	DB	827	U
23	DB	828	U
23	DB	859	G
23	DB	871	U
23	DB	872	U
23	DB	876	C
23	DB	877	A
23	DB	899	A
23	DB	910	A
23	DB	912	C
23	DB	931	U
23	DB	932	U
23	DB	933	A
23	DB	941	A
23	DB	946	C
23	DB	958	U
23	DB	961	C
23	DB	973	A
23	DB	974	G
23	DB	980	A
23	DB	982	C
23	DB	983	A
23	DB	990	A
23	DB	991	C
23	DB	995	C
23	DB	996	A
23	DB	1005	C
23	DB	1012	U
23	DB	1013	C
23	DB	1022	G
23	DB	1023	U
23	DB	1024	G
23	DB	1025	G

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Mol	Chain	Res	Type
23	DB	1033	U
23	DB	1045	C
23	DB	1046	A
23	DB	1047	G
23	DB	1070	A
23	DB	1088	A
23	DB	1090	A
23	DB	1112	G
23	DB	1129	A
23	DB	1132	U
23	DB	1133	A
23	DB	1134	A
23	DB	1135	C
23	DB	1136	G
23	DB	1141	U
23	DB	1157	G
23	DB	1171	G
23	DB	1174	U
23	DB	1175	A
23	DB	1176	U
23	DB	1179	G
23	DB	1195	G
23	DB	1204	A
23	DB	1205	A
23	DB	1206	G
23	DB	1211	C
23	DB	1212	G
23	DB	1237	A
23	DB	1241	A
23	DB	1242	U
23	DB	1248	G
23	DB	1250	G
23	DB	1251	C
23	DB	1253	A
23	DB	1256	G
23	DB	1266	G
23	DB	1271	G
23	DB	1272	A
23	DB	1273	U
23	DB	1275	A
23	DB	1300	G
23	DB	1301	A

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Mol	Chain	Res	Type
23	DB	1318	U
23	DB	1321	A
23	DB	1324	G
23	DB	1325	U
23	DB	1337	G
23	DB	1341	G
23	DB	1365	A
23	DB	1368	G
23	DB	1379	U
23	DB	1383	A
23	DB	1384	A
23	DB	1386	C
23	DB	1396	U
23	DB	1397	U
23	DB	1416	G
23	DB	1419	A
23	DB	1421	G
23	DB	1427	A
23	DB	1428	C
23	DB	1434	A
23	DB	1453	A
23	DB	1454	C
23	DB	1458	U
23	DB	1459	G
23	DB	1460	U
23	DB	1461	C
23	DB	1476	U
23	DB	1477	A
23	DB	1478	G
23	DB	1482	G
23	DB	1490	A
23	DB	1493	C
23	DB	1494	A
23	DB	1504	A
23	DB	1505	A
23	DB	1506	U
23	DB	1507	C
23	DB	1508	A
23	DB	1509	A
23	DB	1523	U
23	DB	1524	G
23	DB	1530	G

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Mol	Chain	Res	Type
23	DB	1532	A
23	DB	1535	A
23	DB	1536	C
23	DB	1552	A
23	DB	1559	U
23	DB	1560	G
23	DB	1567	G
23	DB	1569	A
23	DB	1578	U
23	DB	1583	A
23	DB	1585	C
23	DB	1602	U
23	DB	1603	A
23	DB	1607	C
23	DB	1608	A
23	DB	1610	A
23	DB	1613	G
23	DB	1616	A
23	DB	1618	A
23	DB	1619	G
23	DB	1634	A
23	DB	1635	A
23	DB	1640	A
23	DB	1647	U
23	DB	1648	U
23	DB	1670	C
23	DB	1674	G
23	DB	1681	G
23	DB	1700	A
23	DB	1701	A
23	DB	1715	G
23	DB	1722	A
23	DB	1724	G
23	DB	1727	C
23	DB	1729	U
23	DB	1730	C
23	DB	1731	G
23	DB	1733	G
23	DB	1738	G
23	DB	1756	G
23	DB	1758	U
23	DB	1764	C

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Mol	Chain	Res	Type
23	DB	1773	A
23	DB	1800	C
23	DB	1801	A
23	DB	1808	A
23	DB	1816	C
23	DB	1848	A
23	DB	1869	G
23	DB	1870	C
23	DB	1876	A
23	DB	1884	G
23	DB	1906	G
23	DB	1913	A
23	DB	1914	C
23	DB	1926	U
23	DB	1929	G
23	DB	1930	G
23	DB	1937	A
23	DB	1938	A
23	DB	1939	U
23	DB	1940	U
23	DB	1955	U
23	DB	1964	G
23	DB	1966	A
23	DB	1967	C
23	DB	1970	A
23	DB	1971	U
23	DB	1972	G
23	DB	1991	U
23	DB	1992	G
23	DB	1993	U
23	DB	1996	C
23	DB	1997	C
23	DB	2022	U
23	DB	2023	C
23	DB	2030	A
23	DB	2031	A
23	DB	2032	G
23	DB	2043	C
23	DB	2055	C
23	DB	2056	G
23	DB	2060	A
23	DB	2061	G

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Mol	Chain	Res	Type
23	DB	2062	A
23	DB	2069	G
23	DB	2072	C
23	DB	2093	G
23	DB	2095	A
23	DB	2096	C
23	DB	2100	G
23	DB	2107	G
23	DB	2108	A
23	DB	2109	U
23	DB	2110	G
23	DB	2134	A
23	DB	2136	G
23	DB	2138	G
23	DB	2143	C
23	DB	2144	G
23	DB	2145	C
23	DB	2146	C
23	DB	2147	A
23	DB	2149	U
23	DB	2155	U
23	DB	2156	G
23	DB	2157	G
23	DB	2181	U
23	DB	2188	U
23	DB	2193	G
23	DB	2198	A
23	DB	2199	A
23	DB	2203	U
23	DB	2204	G
23	DB	2210	U
23	DB	2211	A
23	DB	2212	A
23	DB	2213	U
23	DB	2214	C
23	DB	2225	A
23	DB	2238	G
23	DB	2239	G
23	DB	2249	U
23	DB	2250	G
23	DB	2279	G
23	DB	2283	C

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Mol	Chain	Res	Type
23	DB	2287	A
23	DB	2288	A
23	DB	2297	A
23	DB	2304	G
23	DB	2305	U
23	DB	2307	G
23	DB	2308	G
23	DB	2309	A
23	DB	2311	A
23	DB	2320	U
23	DB	2321	U
23	DB	2322	A
23	DB	2324	U
23	DB	2325	G
23	DB	2333	A
23	DB	2336	A
23	DB	2337	G
23	DB	2345	G
23	DB	2347	C
23	DB	2357	G
23	DB	2361	G
23	DB	2372	U
23	DB	2383	G
23	DB	2385	C
23	DB	2396	G
23	DB	2402	U
23	DB	2403	C
23	DB	2406	A
23	DB	2423	U
23	DB	2426	A
23	DB	2427	C
23	DB	2429	G
23	DB	2430	A
23	DB	2431	U
23	DB	2441	U
23	DB	2448	A
23	DB	2472	G
23	DB	2476	A
23	DB	2491	U
23	DB	2492	U
23	DB	2502	G
23	DB	2505	G

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Mol	Chain	Res	Type
23	DB	2506	U
23	DB	2518	A
23	DB	2529	G
23	DB	2543	G
23	DB	2552	U
23	DB	2554	U
23	DB	2566	A
23	DB	2567	G
23	DB	2572	A
23	DB	2573	C
23	DB	2574	G
23	DB	2586	U
23	DB	2602	A
23	DB	2609	U
23	DB	2613	U
23	DB	2629	U
23	DB	2639	A
23	DB	2682	A
23	DB	2689	U
23	DB	2690	U
23	DB	2714	G
23	DB	2726	A
23	DB	2737	G
23	DB	2744	G
23	DB	2748	A
23	DB	2757	A
23	DB	2765	A
23	DB	2778	A
23	DB	2791	G
23	DB	2796	U
23	DB	2798	U
23	DB	2799	A
23	DB	2808	G
23	DB	2820	A
23	DB	2821	A
23	DB	2823	A
23	DB	2835	A
23	DB	2836	U
23	DB	2850	A
23	DB	2859	G
23	DB	2867	G
23	DB	2872	A

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Mol	Chain	Res	Type
23	DB	2873	A
23	DB	2880	C
23	DB	2883	A
23	DB	2894	G
23	DB	2903	U

All (86) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	51	A
1	AA	60	A
1	AA	84	U
1	AA	239	U
1	AA	243	A
1	AA	279	A
1	AA	328	C
1	AA	366	A
1	AA	372	C
1	AA	429	U
1	AA	461	A
1	AA	484	G
1	AA	975	A
1	AA	1049	U
1	AA	1065	U
1	AA	1101	A
1	AA	1201	A
1	AA	1213	A
1	AA	1226	C
1	AA	1302	C
1	AA	1319	A
1	AA	1378	C
1	AA	1397	C
1	AA	1451	U
1	AA	1499	A
22	BA	25	U
23	BB	33	C
23	BB	489	G
23	BB	670	A
23	BB	1210	G
23	BB	1458	U
23	BB	1583	A
23	BB	1847	A

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Mol	Chain	Res	Type
23	BB	1913	A
23	BB	2211	A
23	BB	2282	G
23	BB	2286	G
23	BB	2336	A
23	BB	2425	A
23	BB	2428	G
23	BB	2430	A
23	BB	2756	U
23	BB	2832	U
23	BB	2894	G
1	CA	51	A
1	CA	60	A
1	CA	239	U
1	CA	243	A
1	CA	279	A
1	CA	328	C
1	CA	366	A
1	CA	372	C
1	CA	429	U
1	CA	461	A
1	CA	484	G
1	CA	960	U
1	CA	975	A
1	CA	1049	U
1	CA	1065	U
1	CA	1067	A
1	CA	1201	A
1	CA	1226	C
1	CA	1362	A
1	CA	1397	C
1	CA	1451	U
22	DA	25	U
23	DB	33	C
23	DB	241	A
23	DB	489	G
23	DB	670	A
23	DB	973	A
23	DB	1210	G
23	DB	1320	C
23	DB	1458	U
23	DB	1583	A

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Mol	Chain	Res	Type
23	DB	1847	A
23	DB	2211	A
23	DB	2282	G
23	DB	2286	G
23	DB	2336	A
23	DB	2425	A
23	DB	2428	G
23	DB	2430	A
23	DB	2756	U
23	DB	2832	U
23	DB	2894	G

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 344 ligands modelled in this entry, 342 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
54	SCM	CA	1659	-	23,25,25	1.61	6 (26%)	26,39,39	1.32	2 (7%)
54	SCM	AA	1661	-	23,25,25	1.66	8 (34%)	26,39,39	1.32	2 (7%)

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
54	SCM	CA	1659	-	-	2/4/57/57	0/3/3/3
54	SCM	AA	1661	-	-	2/4/57/57	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
54	CA	1659	SCM	C9-C8	3.10	1.59	1.53
54	AA	1661	SCM	C3-C2	3.03	1.58	1.51
54	AA	1661	SCM	C9-C8	2.91	1.58	1.53
54	CA	1659	SCM	C3-C2	2.82	1.57	1.51
54	AA	1661	SCM	C3-C4	2.56	1.54	1.50
54	AA	1661	SCM	C12-C7	2.36	1.57	1.52
54	CA	1659	SCM	C3-C4	2.35	1.54	1.50
54	CA	1659	SCM	C12-C7	2.28	1.57	1.52
54	AA	1661	SCM	O2B-C12	2.28	1.47	1.44
54	AA	1661	SCM	O5-C5	2.16	1.43	1.39
54	CA	1659	SCM	O2B-C12	2.09	1.47	1.44
54	AA	1661	SCM	C8-N8	2.04	1.51	1.47
54	CA	1659	SCM	C8-N8	2.02	1.51	1.47
54	AA	1661	SCM	C10-N10	-2.02	1.44	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
54	AA	1661	SCM	C1M-N10-C10	-4.88	107.28	114.38
54	CA	1659	SCM	C1M-N10-C10	-4.55	107.76	114.38
54	CA	1659	SCM	C2M-C2-C3	-2.62	108.11	113.22
54	AA	1661	SCM	C2M-C2-C3	-2.42	108.49	113.22

There are no chirality outliers.

All (4) torsion outliers are listed below:

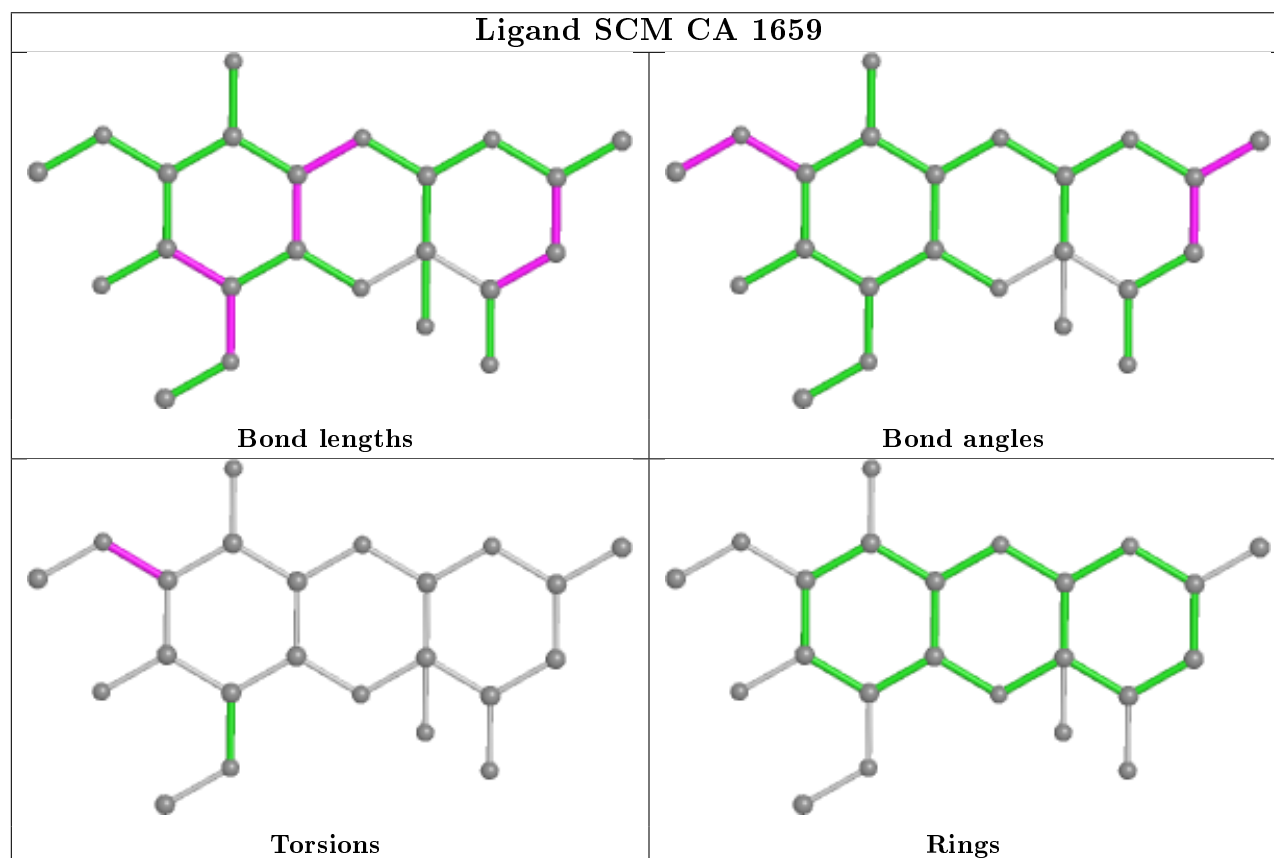
Mol	Chain	Res	Type	Atoms
54	AA	1661	SCM	C9-C10-N10-C1M
54	AA	1661	SCM	C11-C10-N10-C1M
54	CA	1659	SCM	C11-C10-N10-C1M
54	CA	1659	SCM	C9-C10-N10-C1M

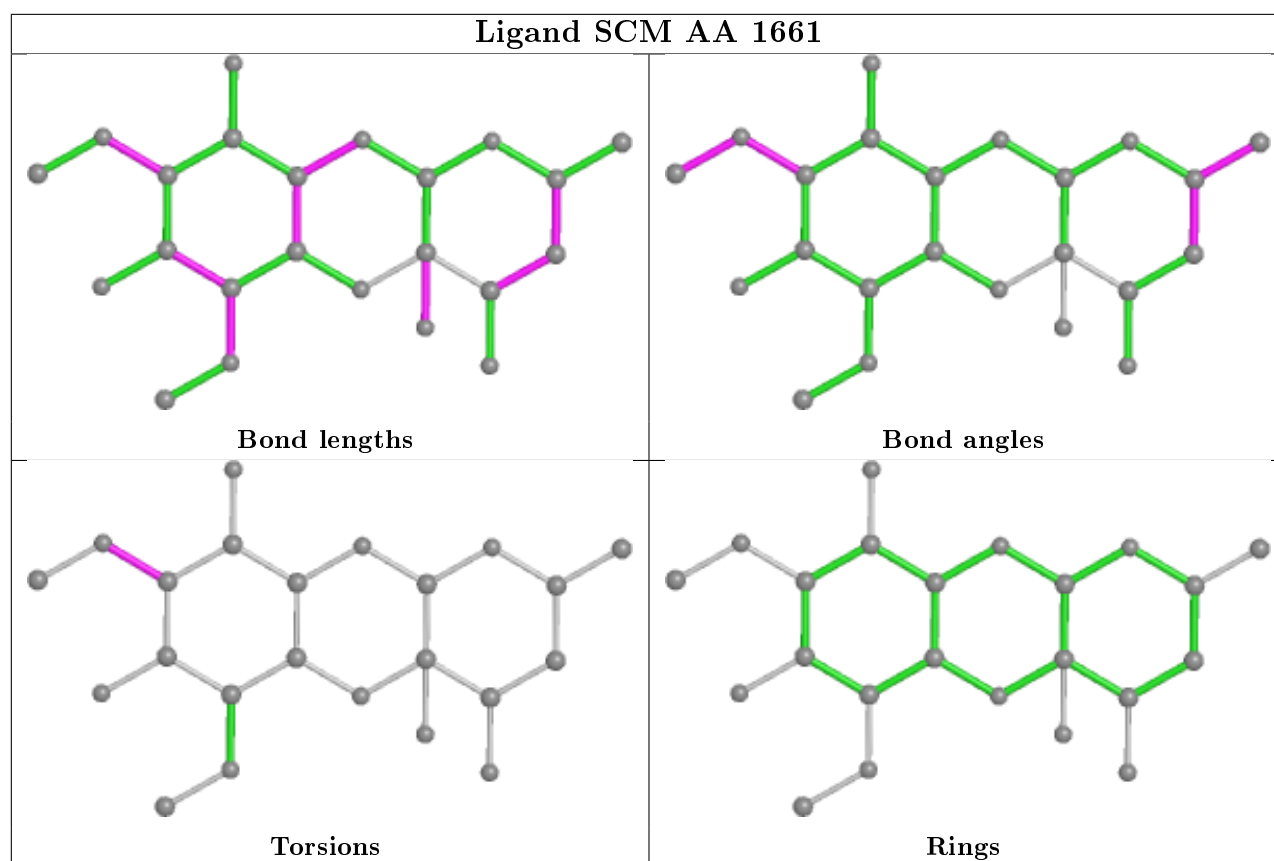
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
54	CA	1659	SCM	1	0
54	AA	1661	SCM	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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	1530/1542 (99%)	-0.77	9 (0%) 89 84	12, 77, 142, 179	0
1	CA	1530/1542 (99%)	-0.76	2 (0%) 95 94	5, 57, 124, 180	0
2	AC	206/232 (88%)	0.22	15 (7%) 15 12	5, 70, 115, 162	0
2	CC	206/232 (88%)	0.53	28 (13%) 3 3	5, 71, 111, 150	0
3	AD	205/205 (100%)	-0.05	12 (5%) 22 18	5, 79, 127, 173	0
3	CD	205/205 (100%)	0.57	20 (9%) 7 7	5, 66, 125, 166	0
4	AE	150/166 (90%)	0.06	4 (2%) 54 44	5, 68, 122, 157	0
4	CE	150/166 (90%)	1.49	51 (34%) 0 0	5, 67, 121, 180	0
5	AF	100/135 (74%)	2.52	62 (62%) 0 0	10, 81, 133, 147	0
5	CF	100/135 (74%)	-0.47	0 100 100	5, 73, 131, 172	0
6	AG	150/178 (84%)	0.74	28 (18%) 1 1	16, 89, 125, 143	0
6	CG	152/178 (85%)	-0.54	0 100 100	6, 79, 125, 172	0
7	AH	129/129 (100%)	0.37	16 (12%) 4 5	19, 77, 130, 158	0
7	CH	129/129 (100%)	0.75	26 (20%) 1 1	5, 62, 116, 158	0
8	AI	127/129 (98%)	0.76	24 (18%) 1 1	6, 83, 117, 155	0
8	CI	127/129 (98%)	0.02	2 (1%) 72 62	5, 80, 122, 157	0
9	AJ	98/103 (95%)	0.49	4 (4%) 37 30	9, 79, 126, 147	0
9	CJ	98/103 (95%)	1.28	29 (29%) 0 0	16, 82, 113, 137	0
10	AK	117/128 (91%)	0.23	4 (3%) 45 36	5, 72, 117, 155	0
10	CK	117/128 (91%)	-0.54	0 100 100	5, 67, 123, 136	0
11	AL	123/123 (100%)	1.01	30 (24%) 0 0	6, 79, 123, 158	0
11	CL	123/123 (100%)	0.26	6 (4%) 29 25	5, 54, 103, 151	0
12	AM	114/117 (97%)	0.95	22 (19%) 1 1	23, 96, 137, 169	0
12	CM	113/117 (96%)	-0.21	5 (4%) 34 29	22, 96, 142, 162	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AP	82/82 (100%)	1.48	25 (30%) 0 0	15, 82, 130, 163	0
13	CP	80/82 (97%)	0.05	5 (6%) 20 16	5, 58, 126, 180	0
14	AQ	80/83 (96%)	1.25	23 (28%) 0 0	15, 92, 145, 154	0
14	CQ	81/83 (97%)	0.26	3 (3%) 41 33	5, 72, 119, 149	0
15	AR	55/74 (74%)	0.81	8 (14%) 2 3	5, 73, 129, 164	0
15	CR	55/74 (74%)	0.23	3 (5%) 25 22	5, 66, 127, 143	0
16	AS	79/91 (86%)	1.98	39 (49%) 0 0	52, 98, 142, 167	0
16	CS	80/91 (87%)	0.50	13 (16%) 1 2	41, 95, 133, 159	0
17	AT	85/86 (98%)	-0.68	0 100 100	19, 88, 126, 156	0
17	CT	85/86 (98%)	-0.28	1 (1%) 79 70	5, 66, 113, 154	0
18	AB	218/240 (90%)	-0.07	10 (4%) 32 27	12, 81, 120, 160	0
18	CB	218/240 (90%)	0.92	50 (22%) 0 0	5, 87, 133, 163	0
19	AU	51/70 (72%)	0.78	8 (15%) 2 2	20, 94, 139, 153	0
19	CU	51/70 (72%)	0.11	2 (3%) 39 31	57, 96, 137, 171	0
20	AO	88/89 (98%)	1.00	11 (12%) 3 5	5, 73, 118, 177	0
20	CO	88/89 (98%)	-0.48	0 100 100	5, 56, 108, 135	0
21	AN	96/100 (96%)	1.33	30 (31%) 0 0	5, 84, 128, 155	0
21	CN	96/100 (96%)	1.13	24 (25%) 0 0	5, 75, 129, 145	0
22	BA	117/120 (97%)	0.07	2 (1%) 70 60	35, 77, 125, 167	0
22	DA	117/120 (97%)	-0.56	1 (0%) 84 77	31, 87, 133, 176	0
23	BB	2841/2904 (97%)	-0.46	23 (0%) 86 79	5, 60, 136, 180	0
23	DB	2841/2904 (97%)	-0.52	23 (0%) 86 79	5, 51, 136, 180	0
24	BI	141/141 (100%)	2.65	68 (48%) 0 0	60, 135, 178, 180	0
24	DI	141/141 (100%)	1.29	38 (26%) 0 0	66, 135, 180, 180	0
25	BC	271/272 (99%)	1.41	88 (32%) 0 0	5, 61, 109, 132	0
25	DC	271/272 (99%)	0.80	54 (19%) 1 1	5, 45, 99, 144	0
26	BD	209/209 (100%)	0.49	29 (13%) 2 3	5, 71, 118, 148	0
26	DD	209/209 (100%)	0.91	38 (18%) 1 2	5, 60, 110, 168	0
27	BK	121/123 (98%)	2.08	65 (53%) 0 0	5, 75, 125, 159	0
27	DK	121/123 (98%)	1.05	27 (22%) 0 1	5, 45, 112, 150	0
28	BP	114/114 (100%)	1.54	47 (41%) 0 0	7, 82, 125, 155	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	DP	114/114 (100%)	0.40	6 (5%) 26 23	5, 64, 118, 139	0
29	BE	201/201 (100%)	0.81	33 (16%) 1 2	5, 65, 128, 164	0
29	DE	201/201 (100%)	1.23	58 (28%) 0 0	5, 73, 123, 160	0
30	BY	58/58 (100%)	-0.26	1 (1%) 70 60	5, 75, 129, 160	0
30	DY	58/58 (100%)	-0.23	3 (5%) 27 24	5, 72, 116, 147	0
31	B0	56/56 (100%)	0.04	2 (3%) 42 34	5, 78, 118, 147	0
31	D0	56/56 (100%)	0.18	1 (1%) 68 60	5, 62, 129, 153	0
32	B4	38/38 (100%)	0.95	8 (21%) 1 1	27, 81, 131, 145	0
32	D4	38/38 (100%)	-0.34	0 100 100	5, 67, 106, 117	0
33	B1	50/54 (92%)	0.62	4 (8%) 12 11	17, 70, 120, 134	0
33	D1	50/54 (92%)	0.34	5 (10%) 7 7	17, 73, 116, 137	0
34	B3	64/64 (100%)	-0.23	1 (1%) 72 62	5, 68, 103, 129	0
34	D3	64/64 (100%)	1.36	22 (34%) 0 0	5, 55, 86, 122	0
35	BV	94/94 (100%)	0.64	14 (14%) 2 3	5, 81, 126, 152	0
35	DV	94/94 (100%)	0.94	21 (22%) 0 1	5, 88, 120, 160	0
36	B2	46/46 (100%)	0.77	5 (10%) 5 5	5, 53, 104, 141	0
36	D2	46/46 (100%)	1.64	20 (43%) 0 0	9, 48, 110, 141	0
37	BL	143/144 (99%)	-0.16	2 (1%) 75 66	5, 67, 121, 145	0
37	DL	143/144 (99%)	1.63	57 (39%) 0 0	5, 63, 111, 145	0
38	BM	136/136 (100%)	0.91	21 (15%) 2 2	7, 68, 121, 179	0
38	DM	136/136 (100%)	1.31	39 (28%) 0 0	5, 65, 117, 144	0
39	BX	63/63 (100%)	1.00	13 (20%) 1 1	16, 79, 128, 159	0
39	DX	63/63 (100%)	1.09	9 (14%) 2 3	16, 91, 144, 169	0
40	BH	149/149 (100%)	3.05	87 (58%) 0 0	7, 104, 149, 180	0
40	DH	149/149 (100%)	0.85	26 (17%) 1 2	5, 91, 131, 162	0
41	BJ	142/142 (100%)	0.94	29 (20%) 1 1	5, 77, 118, 140	0
41	DJ	142/142 (100%)	0.27	10 (7%) 16 13	5, 70, 119, 173	0
42	BN	120/127 (94%)	0.84	17 (14%) 2 3	5, 69, 111, 154	0
42	DN	120/127 (94%)	0.02	6 (5%) 28 25	5, 51, 94, 163	0
43	BO	116/117 (99%)	-0.08	7 (6%) 21 17	6, 80, 109, 172	0
43	DO	116/117 (99%)	0.90	20 (17%) 1 2	5, 83, 136, 158	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	BQ	117/117 (100%)	-0.62	1 (0%) 84 77	5, 62, 115, 144	0
44	DQ	117/117 (100%)	0.46	12 (10%) 6 6	5, 60, 111, 154	0
45	BS	110/110 (100%)	0.39	7 (6%) 19 15	5, 62, 121, 148	0
45	DS	110/110 (100%)	1.42	36 (32%) 0 0	5, 64, 127, 156	0
46	BU	102/103 (99%)	0.84	16 (15%) 2 2	12, 80, 125, 157	0
46	DU	102/103 (99%)	-0.41	0 100 100	8, 94, 127, 149	0
47	BF	178/178 (100%)	0.73	31 (17%) 1 2	29, 100, 146, 180	0
47	DF	178/178 (100%)	2.23	89 (50%) 0 0	12, 93, 142, 163	0
48	BG	176/176 (100%)	0.76	31 (17%) 1 2	18, 94, 133, 171	0
48	DG	176/176 (100%)	0.40	24 (13%) 3 3	8, 90, 136, 166	0
49	BR	103/103 (100%)	-0.03	4 (3%) 39 31	5, 83, 123, 133	0
49	DR	103/103 (100%)	0.13	5 (4%) 29 25	10, 76, 135, 149	0
50	BT	93/100 (93%)	-0.07	2 (2%) 62 52	8, 83, 130, 165	0
50	DT	93/100 (93%)	2.27	51 (54%) 0 0	5, 84, 144, 172	0
51	BZ	77/78 (98%)	0.57	12 (15%) 2 2	5, 63, 120, 142	0
51	DZ	77/78 (98%)	0.01	1 (1%) 77 68	5, 56, 106, 120	0
52	BW	79/84 (94%)	0.45	6 (7%) 13 12	5, 75, 124, 180	0
52	DW	79/84 (94%)	0.69	9 (11%) 5 5	5, 71, 121, 166	0
All	All	20417/21046 (97%)	0.14	2041 (9%) 7 7	5, 70, 134, 180	0

All (2041) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
24	BI	17	ALA	14.4
23	BB	546	U	12.7
24	BI	18	ASN	12.4
40	BH	124	THR	12.3
24	BI	51	GLY	12.2
23	BB	1731	G	12.2
23	BB	1730	C	11.7
24	BI	49	GLU	11.5
40	BH	142	VAL	10.3
23	BB	140	C	10.3
40	BH	93	SER	10.0
24	BI	52	LEU	9.7
40	BH	130	VAL	9.4

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Mol	Chain	Res	Type	RSRZ
13	AP	82	ALA	9.3
40	BH	94	ILE	8.9
33	D1	52	LYS	8.7
15	AR	19	GLU	8.6
40	BH	131	SER	8.6
24	BI	48	ILE	8.4
40	BH	144	VAL	8.4
24	BI	14	ALA	8.3
39	DX	63	ALA	8.3
39	DX	62	GLY	8.1
40	BH	120	GLY	8.0
40	BH	80	ILE	8.0
3	CD	163	GLN	7.9
8	AI	129	ARG	7.8
24	BI	54	ILE	7.8
16	AS	40	PHE	7.8
13	AP	81	ALA	7.7
52	DW	84	GLU	7.7
50	DT	92	ASN	7.6
24	BI	34	ILE	7.6
50	DT	5	GLU	7.5
23	BB	1728	C	7.5
40	BH	128	HIS	7.5
16	AS	41	PRO	7.5
40	BH	125	THR	7.5
33	B1	52	LYS	7.4
24	BI	53	PRO	7.4
24	DI	95	ASP	7.3
47	DF	44	ALA	7.3
50	DT	72	GLN	7.2
47	DF	173	ASP	7.2
21	CN	30	ILE	7.0
40	BH	117	LEU	7.0
24	BI	19	PRO	7.0
40	BH	149	GLU	7.0
24	BI	16	MET	6.9
13	AP	80	LYS	6.9
40	BH	133	GLN	6.9
24	BI	50	LYS	6.8
24	BI	15	GLY	6.7
23	BB	1727	C	6.7
39	BX	62	GLY	6.7

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Mol	Chain	Res	Type	RSRZ
47	DF	10	GLU	6.7
51	BZ	78	TYR	6.7
40	BH	146	VAL	6.7
24	BI	20	SER	6.6
50	DT	91	GLN	6.6
34	D3	19	GLY	6.5
27	BK	122	VAL	6.5
47	DF	139	GLU	6.5
24	BI	21	PRO	6.5
40	BH	102	ALA	6.5
47	BF	127	TYR	6.5
24	BI	11	GLN	6.5
47	DF	174	PHE	6.4
24	DI	99	LYS	6.4
29	DE	124	PHE	6.4
24	DI	98	GLY	6.4
40	BH	145	ASN	6.4
24	BI	58	ILE	6.3
29	BE	155	GLU	6.3
12	AM	114	PRO	6.3
40	BH	86	ASP	6.3
5	AF	90	MET	6.3
18	CB	161	PHE	6.3
41	BJ	63	ALA	6.2
40	BH	87	GLU	6.2
40	BH	121	VAL	6.2
40	BH	123	ARG	6.2
5	AF	8	PHE	6.2
40	BH	115	VAL	6.1
40	BH	92	GLY	6.1
5	AF	66	ALA	6.1
25	BC	114	GLN	6.1
40	BH	126	GLY	6.0
40	BH	119	ASN	6.0
40	BH	88	GLY	6.0
40	BH	45	GLU	6.0
40	BH	103	VAL	5.9
6	AG	79	VAL	5.9
38	DM	1	MET	5.9
50	DT	53	VAL	5.9
24	BI	13	ALA	5.9
28	BP	71	ARG	5.9

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Mol	Chain	Res	Type	RSRZ
16	CS	26	ASP	5.8
24	BI	68	PHE	5.8
47	DF	127	TYR	5.8
4	CE	146	MET	5.8
7	CH	98	LEU	5.8
52	DW	18	LYS	5.8
21	CN	20	PHE	5.8
11	AL	85	ARG	5.8
23	BB	139	U	5.7
21	CN	42	ASN	5.7
22	DA	88	C	5.7
40	BH	116	ARG	5.7
48	DG	40	VAL	5.7
23	DB	1730	C	5.7
24	DI	138	VAL	5.7
26	DD	32	ASN	5.7
18	CB	163	ILE	5.7
50	DT	11	LEU	5.7
24	DI	137	LEU	5.6
47	DF	55	ASP	5.6
38	DM	105	MET	5.6
21	CN	31	SER	5.6
24	BI	59	THR	5.6
47	DF	19	PHE	5.6
5	AF	88	MET	5.6
38	DM	60	GLN	5.6
5	AF	5	GLU	5.5
47	DF	171	ALA	5.5
52	BW	83	ALA	5.5
52	DW	19	ARG	5.5
29	DE	155	GLU	5.5
21	CN	23	ARG	5.4
38	DM	32	GLY	5.4
40	DH	82	SER	5.4
8	CI	129	ARG	5.4
13	AP	52	LEU	5.4
12	AM	79	LEU	5.4
50	DT	12	ARG	5.4
37	DL	92	LEU	5.4
11	CL	123	ALA	5.4
24	BI	25	PRO	5.4
23	BB	1729	U	5.4

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Mol	Chain	Res	Type	RSRZ
23	BB	715	A	5.4
24	DI	84	GLY	5.4
50	DT	70	HIS	5.4
24	BI	97	VAL	5.4
4	CE	81	GLN	5.4
8	AI	56	MET	5.3
26	BD	131	ASP	5.3
38	BM	103	TYR	5.3
9	CJ	101	SER	5.3
23	BB	1726	C	5.3
27	BK	110	GLU	5.3
8	AI	51	LEU	5.3
19	AU	3	ILE	5.3
11	CL	13	ARG	5.3
4	CE	85	LYS	5.3
9	CJ	102	LEU	5.3
26	BD	146	ILE	5.3
21	CN	19	TYR	5.3
6	AG	85	GLN	5.3
37	DL	45	GLY	5.2
48	DG	52	GLY	5.2
32	B4	7	VAL	5.2
24	BI	85	ILE	5.2
27	BK	121	GLU	5.2
7	CH	129	ALA	5.2
37	DL	91	ASP	5.2
21	CN	34	ASN	5.2
24	BI	37	PHE	5.2
27	BK	103	VAL	5.2
40	BH	141	LYS	5.2
9	CJ	35	GLN	5.1
7	AH	129	ALA	5.1
28	BP	61	ARG	5.1
23	BB	1067	A	5.1
5	AF	63	ASN	5.1
47	DF	43	ILE	5.1
40	BH	140	ALA	5.1
29	BE	124	PHE	5.1
37	DL	58	TYR	5.1
5	AF	67	PRO	5.1
24	BI	12	VAL	5.1
18	CB	56	LEU	5.1

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Mol	Chain	Res	Type	RSRZ
35	BV	74	ALA	5.0
40	BH	85	GLY	5.0
24	BI	45	THR	5.0
4	CE	12	GLU	5.0
13	CP	47	GLU	5.0
15	AR	22	TYR	5.0
47	DF	157	THR	5.0
16	AS	43	MET	5.0
34	D3	13	PHE	5.0
41	BJ	64	VAL	5.0
47	DF	30	VAL	5.0
37	DL	108	ALA	5.0
47	DF	172	PHE	5.0
3	CD	22	SER	4.9
40	BH	132	PHE	4.9
47	DF	140	ILE	4.9
25	DC	81	GLU	4.9
40	BH	118	PRO	4.9
24	DI	136	GLY	4.9
29	BE	14	VAL	4.9
43	DO	24	THR	4.9
40	BH	84	ALA	4.9
4	CE	36	THR	4.9
24	BI	27	LEU	4.9
27	BK	102	PRO	4.9
11	AL	13	ARG	4.9
39	DX	22	LEU	4.9
18	CB	68	PHE	4.9
36	D2	1	MET	4.8
33	B1	51	ALA	4.8
47	DF	42	ALA	4.8
34	D3	51	LYS	4.8
40	BH	122	LEU	4.8
15	CR	19	GLU	4.8
4	CE	158	LYS	4.8
38	BM	32	GLY	4.8
5	AF	35	LYS	4.8
16	AS	73	PHE	4.8
41	DJ	64	VAL	4.8
40	BH	101	ASP	4.8
40	BH	82	SER	4.8
50	DT	73	ARG	4.8

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Mol	Chain	Res	Type	RSRZ
40	BH	98	ASP	4.8
5	AF	7	VAL	4.7
24	BI	26	ALA	4.7
21	AN	28	ALA	4.7
11	AL	14	LYS	4.7
19	AU	35	GLU	4.7
6	AG	80	GLY	4.7
26	DD	2	ILE	4.7
29	BE	119	ILE	4.7
2	CC	165	GLU	4.7
9	CJ	81	GLU	4.7
25	BC	1	ALA	4.7
27	BK	104	THR	4.7
24	BI	29	GLN	4.7
51	DZ	78	TYR	4.7
50	DT	71	GLY	4.7
29	BE	153	LEU	4.7
38	DM	2	LEU	4.7
25	BC	123	ILE	4.7
47	DF	82	TYR	4.7
25	BC	22	GLU	4.7
40	BH	60	GLU	4.7
47	BF	139	GLU	4.7
47	DF	164	GLU	4.6
37	DL	123	ARG	4.6
8	AI	49	GLN	4.6
40	BH	148	ALA	4.6
5	AF	9	MET	4.6
52	BW	84	GLU	4.6
28	BP	47	ILE	4.6
37	DL	49	GLY	4.6
38	DM	30	SER	4.6
9	CJ	79	PRO	4.6
47	DF	27	VAL	4.6
4	CE	114	LEU	4.6
11	AL	11	ARG	4.6
49	BR	46	GLU	4.6
28	BP	42	PHE	4.6
24	DI	85	ILE	4.5
5	AF	70	VAL	4.5
40	DH	1	MET	4.5
18	CB	160	LEU	4.5

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Mol	Chain	Res	Type	RSRZ
21	CN	18	LYS	4.5
47	DF	11	VAL	4.5
48	DG	55	ASP	4.5
41	BJ	86	GLN	4.5
11	AL	19	ASN	4.5
42	BN	83	LEU	4.5
38	DM	103	TYR	4.5
4	CE	80	LEU	4.5
26	DD	1	MET	4.5
47	DF	33	ILE	4.5
45	DS	12	SER	4.5
21	CN	22	LYS	4.5
6	AG	8	GLN	4.5
29	BE	188	MET	4.5
18	CB	185	ILE	4.5
40	DH	149	GLU	4.5
35	BV	56	PHE	4.5
37	DL	57	LEU	4.5
5	AF	10	VAL	4.4
47	DF	28	PRO	4.4
6	AG	20	GLU	4.4
28	BP	70	GLU	4.4
27	BK	8	LEU	4.4
4	CE	117	ALA	4.4
21	CN	17	ASP	4.4
40	BH	81	ALA	4.4
21	AN	23	ARG	4.4
27	BK	84	CYS	4.4
38	BM	33	LEU	4.4
42	BN	21	PHE	4.4
51	BZ	76	GLU	4.4
26	DD	95	SER	4.4
40	DH	12	LEU	4.4
48	BG	42	VAL	4.4
51	BZ	77	LYS	4.4
31	D0	56	LYS	4.4
29	DE	169	VAL	4.4
44	DQ	4	LYS	4.4
48	DG	41	GLU	4.3
47	DF	41	GLU	4.3
47	DF	75	GLY	4.3
45	DS	104	THR	4.3

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Mol	Chain	Res	Type	RSRZ
12	AM	55	LEU	4.3
5	AF	6	ILE	4.3
23	DB	544	C	4.3
25	BC	113	ASP	4.3
2	AC	45	GLU	4.3
9	CJ	34	ALA	4.3
21	AN	24	ALA	4.3
26	DD	47	ALA	4.3
40	BH	95	GLY	4.3
24	BI	3	LYS	4.3
37	DL	82	LEU	4.3
9	CJ	36	VAL	4.3
25	DC	57	HIS	4.3
47	DF	7	TYR	4.3
47	DF	175	PRO	4.3
14	AQ	45	VAL	4.3
50	DT	10	VAL	4.3
13	AP	7	ALA	4.3
52	DW	83	ALA	4.3
21	AN	20	PHE	4.3
24	BI	47	SER	4.3
27	DK	18	ARG	4.3
47	DF	24	VAL	4.3
23	BB	1537	G	4.3
29	BE	150	THR	4.3
50	DT	87	LEU	4.3
28	BP	48	ALA	4.3
52	DW	17	ALA	4.3
12	AM	112	ARG	4.2
18	CB	184	ALA	4.2
50	DT	35	ALA	4.2
40	BH	99	ILE	4.2
33	D1	26	LYS	4.2
24	DI	97	VAL	4.2
39	BX	8	GLU	4.2
24	BI	38	CYS	4.2
45	DS	81	SER	4.2
40	DH	18	GLN	4.2
50	DT	43	ILE	4.2
3	AD	106	PHE	4.2
48	DG	114	HIS	4.2
18	CB	212	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
47	DF	124	ARG	4.2
47	DF	169	LEU	4.2
38	DM	106	ASP	4.2
45	DS	93	ALA	4.2
33	D1	27	ARG	4.2
48	BG	147	LEU	4.2
38	DM	29	GLY	4.2
25	DC	1	ALA	4.2
6	AG	61	PHE	4.2
16	AS	29	PRO	4.2
25	BC	265	PHE	4.2
37	DL	141	LYS	4.2
29	DE	60	TRP	4.2
5	AF	65	GLU	4.2
14	AQ	8	GLN	4.2
41	DJ	44	TYR	4.2
40	BH	129	GLU	4.2
24	BI	69	VAL	4.2
50	DT	15	HIS	4.2
11	AL	47	ALA	4.1
12	AM	82	LEU	4.1
5	AF	1	MET	4.1
26	BD	186	LEU	4.1
11	AL	1	ALA	4.1
5	AF	4	TYR	4.1
29	DE	148	ILE	4.1
50	DT	90	GLY	4.1
24	BI	70	THR	4.1
28	BP	43	GLU	4.1
24	DI	89	SER	4.1
45	DS	103	ILE	4.1
28	BP	99	LEU	4.1
48	BG	40	VAL	4.1
6	AG	10	LYS	4.1
22	BA	118	C	4.1
37	DL	51	GLU	4.1
26	DD	87	GLY	4.1
5	AF	34	GLY	4.1
24	BI	32	VAL	4.1
24	BI	46	ASP	4.1
38	DM	134	THR	4.1
50	DT	93	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
28	BP	73	PHE	4.1
3	CD	24	VAL	4.1
35	BV	57	TYR	4.1
45	DS	94	ASP	4.1
38	DM	64	TRP	4.1
47	DF	12	VAL	4.1
47	DF	134	GLN	4.0
18	CB	200	PRO	4.0
26	DD	165	MET	4.0
5	AF	87	SER	4.0
50	DT	81	LYS	4.0
5	AF	62	MET	4.0
26	BD	25	THR	4.0
4	CE	13	LYS	4.0
38	BM	105	MET	4.0
38	BM	1	MET	4.0
48	BG	120	ILE	4.0
48	BG	176	LYS	4.0
24	BI	141	ASP	4.0
26	DD	52	THR	4.0
34	D3	21	PHE	4.0
24	BI	22	PRO	4.0
40	BH	127	GLU	4.0
25	BC	116	GLN	4.0
25	BC	182	LYS	4.0
50	DT	36	LYS	4.0
45	DS	92	ARG	4.0
40	DH	81	ALA	4.0
4	CE	11	GLN	4.0
25	DC	59	GLN	4.0
40	BH	79	THR	4.0
26	DD	96	ILE	4.0
29	BE	10	SER	4.0
47	DF	90	LEU	4.0
29	DE	65	THR	4.0
45	BS	74	ILE	4.0
12	AM	113	LYS	4.0
3	CD	164	ARG	4.0
5	AF	54	LEU	4.0
25	DC	5	CYS	4.0
40	BH	90	LEU	4.0
50	DT	14	PRO	4.0

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Mol	Chain	Res	Type	RSRZ
6	AG	84	TYR	4.0
18	CB	195	VAL	3.9
38	DM	37	GLY	3.9
27	BK	73	ASP	3.9
39	BX	63	ALA	3.9
24	DI	93	ASN	3.9
28	BP	67	GLU	3.9
16	AS	39	ILE	3.9
37	DL	86	GLU	3.9
48	DG	37	ASN	3.9
24	DI	94	LYS	3.9
7	CH	71	VAL	3.9
24	BI	55	PRO	3.9
29	DE	201	ALA	3.9
1	AA	1362	A	3.9
5	AF	64	VAL	3.9
11	AL	61	GLU	3.9
12	CM	44	ILE	3.9
40	BH	1	MET	3.9
25	DC	62	ARG	3.9
4	CE	10	LEU	3.9
47	DF	58	ALA	3.9
37	DL	59	ARG	3.9
40	BH	56	ALA	3.9
29	DE	149	ILE	3.9
29	BE	1	MET	3.9
31	B0	54	ILE	3.8
38	DM	104	GLU	3.8
16	AS	11	ASP	3.8
38	DM	31	PHE	3.8
4	CE	147	ASN	3.8
16	CS	20	LYS	3.8
8	AI	57	VAL	3.8
27	BK	77	ILE	3.8
4	CE	148	SER	3.8
24	BI	35	MET	3.8
39	BX	5	GLU	3.8
8	AI	47	VAL	3.8
18	CB	69	VAL	3.8
48	DG	57	TYR	3.8
40	BH	57	LYS	3.8
25	BC	20	ASN	3.8

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Mol	Chain	Res	Type	RSRZ
46	BU	71	ILE	3.8
24	DI	96	LYS	3.8
29	DE	88	ARG	3.8
24	BI	87	SER	3.8
29	DE	55	SER	3.8
7	CH	99	GLY	3.8
20	AO	43	PHE	3.8
40	BH	46	PHE	3.8
18	CB	164	ASP	3.8
24	DI	83	ALA	3.8
24	DI	132	ALA	3.8
25	BC	115	ILE	3.8
5	AF	11	HIS	3.8
5	AF	42	TRP	3.8
6	AG	15	PRO	3.8
29	DE	70	SER	3.8
37	DL	118	THR	3.8
38	BM	136	MET	3.8
24	BI	6	ALA	3.8
6	AG	58	LEU	3.8
16	AS	59	VAL	3.8
50	DT	13	ALA	3.8
50	DT	34	VAL	3.8
5	AF	94	HIS	3.8
8	AI	128	LYS	3.8
16	AS	77	ARG	3.8
25	BC	17	LYS	3.7
5	AF	78	PHE	3.7
45	DS	11	ARG	3.7
25	DC	22	GLU	3.7
14	AQ	20	ILE	3.7
27	BK	94	PRO	3.7
29	BE	118	LEU	3.7
47	DF	149	ARG	3.7
19	CU	23	GLU	3.7
25	DC	33	LEU	3.7
8	AI	50	PRO	3.7
16	AS	42	ASN	3.7
36	D2	37	LYS	3.7
18	CB	198	VAL	3.7
25	DC	79	ARG	3.7
37	DL	50	PHE	3.7

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Mol	Chain	Res	Type	RSRZ
26	DD	75	ALA	3.7
4	CE	115	GLU	3.7
50	DT	31	VAL	3.7
21	AN	29	ILE	3.7
29	BE	125	SER	3.7
47	DF	155	ILE	3.7
5	AF	93	LYS	3.7
41	BJ	24	THR	3.7
11	AL	86	VAL	3.7
14	AQ	58	VAL	3.7
4	CE	95	MET	3.7
25	BC	269	ARG	3.7
37	DL	47	ARG	3.7
5	AF	100	SER	3.7
45	DS	5	ALA	3.7
5	AF	25	TYR	3.7
5	AF	91	ARG	3.7
26	DD	93	GLY	3.7
27	BK	80	ASP	3.7
24	BI	24	GLY	3.7
40	BH	91	PHE	3.7
16	AS	60	PHE	3.7
36	D2	29	GLN	3.7
5	AF	68	GLN	3.7
25	DC	34	GLU	3.7
23	DB	139	U	3.7
11	AL	63	THR	3.7
6	AG	11	ILE	3.6
41	BJ	87	ALA	3.6
43	DO	3	LYS	3.6
11	AL	93	ARG	3.6
4	CE	43	GLY	3.6
24	BI	28	GLY	3.6
26	BD	188	LEU	3.6
7	AH	126	CYS	3.6
48	DG	88	LEU	3.6
46	BU	83	GLY	3.6
29	DE	74	LYS	3.6
38	DM	28	PHE	3.6
41	BJ	54	ILE	3.6
47	DF	86	CYS	3.6
36	D2	2	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
25	BC	181	ARG	3.6
27	BK	98	ARG	3.6
48	DG	176	LYS	3.6
27	BK	64	ARG	3.6
14	AQ	53	GLY	3.6
4	CE	144	GLU	3.6
47	DF	89	THR	3.6
29	BE	143	LEU	3.6
43	DO	2	ASP	3.6
4	CE	74	ALA	3.6
25	DC	213	ARG	3.6
27	BK	46	ALA	3.6
38	DM	61	GLY	3.6
47	BF	75	GLY	3.6
28	BP	75	THR	3.6
42	BN	70	THR	3.6
46	BU	93	ARG	3.6
7	AH	74	ILE	3.6
24	BI	60	VAL	3.6
40	BH	55	GLU	3.6
50	DT	51	PHE	3.6
41	BJ	142	ILE	3.6
9	CJ	26	VAL	3.6
40	BH	147	VAL	3.6
37	DL	61	LEU	3.6
46	BU	84	PHE	3.6
16	AS	28	LYS	3.6
29	DE	188	MET	3.6
47	DF	160	LYS	3.6
12	AM	85	TYR	3.6
52	BW	64	GLY	3.6
50	DT	6	ARG	3.6
12	AM	6	ILE	3.6
27	BK	35	VAL	3.6
35	DV	74	ALA	3.6
16	AS	65	MET	3.5
29	BE	120	VAL	3.5
29	DE	72	SER	3.5
7	CH	127	TYR	3.5
29	DE	153	LEU	3.5
47	DF	18	GLU	3.5
28	BP	41	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
37	DL	55	MET	3.5
25	DC	4	LYS	3.5
26	DD	33	ARG	3.5
29	DE	61	ARG	3.5
4	AE	94	PHE	3.5
5	AF	83	ALA	3.5
16	AS	79	TYR	3.5
29	DE	75	SER	3.5
37	DL	76	GLU	3.5
3	CD	16	THR	3.5
37	DL	77	ILE	3.5
6	AG	12	LEU	3.5
36	D2	3	ARG	3.5
16	AS	38	THR	3.5
41	DJ	45	THR	3.5
18	AB	195	VAL	3.5
50	DT	30	ILE	3.5
3	AD	173	ASP	3.5
42	DN	83	LEU	3.5
42	DN	73	ASN	3.5
37	DL	85	VAL	3.5
29	DE	134	LEU	3.5
36	B2	1	MET	3.5
46	BU	59	GLU	3.5
6	AG	76	SER	3.5
37	DL	101	ILE	3.5
2	CC	133	MET	3.5
26	DD	35	THR	3.5
26	BD	111	GLY	3.5
37	BL	144	GLU	3.5
47	DF	156	THR	3.5
39	BX	13	GLU	3.5
35	DV	51	GLN	3.5
38	BM	102	LEU	3.5
47	DF	87	LYS	3.5
41	DJ	1	MET	3.5
42	BN	24	MET	3.5
8	AI	38	PHE	3.5
15	AR	23	LYS	3.5
40	BH	104	THR	3.5
40	BH	134	VAL	3.5
38	DM	33	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
45	BS	103	ILE	3.5
4	CE	79	THR	3.5
41	DJ	87	ALA	3.5
36	D2	46	LYS	3.5
28	BP	19	PHE	3.5
28	BP	46	VAL	3.5
29	DE	187	VAL	3.5
33	D1	6	GLU	3.5
18	CB	74	ALA	3.4
27	BK	15	GLY	3.4
47	DF	103	ILE	3.4
25	BC	93	VAL	3.4
35	BV	91	PHE	3.4
13	CP	52	LEU	3.4
50	DT	33	LYS	3.4
50	DT	44	LYS	3.4
43	DO	87	ILE	3.4
4	CE	37	VAL	3.4
24	BI	40	ALA	3.4
21	CN	26	LEU	3.4
24	BI	33	ASN	3.4
26	BD	187	LEU	3.4
29	DE	177	PRO	3.4
35	BV	84	PRO	3.4
18	CB	199	ILE	3.4
26	DD	48	ILE	3.4
34	D3	15	LYS	3.4
50	DT	82	LYS	3.4
4	CE	73	VAL	3.4
25	BC	131	MET	3.4
26	DD	3	GLY	3.4
16	AS	2	ARG	3.4
18	CB	183	PHE	3.4
36	D2	33	ARG	3.4
21	AN	40	ARG	3.4
2	CC	167	TYR	3.4
25	BC	193	GLU	3.4
40	BH	12	LEU	3.4
36	D2	26	ASN	3.4
41	BJ	102	GLU	3.4
24	DI	52	LEU	3.4
40	DH	122	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
43	DO	88	LYS	3.4
28	BP	58	PHE	3.4
21	AN	45	LEU	3.4
19	CU	3	ILE	3.4
24	DI	47	SER	3.4
9	CJ	85	ASP	3.4
47	BF	112	ASP	3.4
37	DL	117	THR	3.4
3	CD	108	ALA	3.4
40	BH	100	ALA	3.4
3	AD	178	GLU	3.4
40	BH	76	GLU	3.4
48	BG	161	VAL	3.4
8	AI	127	SER	3.4
12	CM	1	ALA	3.4
29	BE	148	ILE	3.4
9	CJ	84	VAL	3.4
35	DV	91	PHE	3.4
45	DS	100	THR	3.4
11	CL	14	LYS	3.3
18	CB	192	PRO	3.3
25	BC	122	ALA	3.3
25	DC	228	ASP	3.3
47	DF	21	TYR	3.3
2	CC	198	LYS	3.3
25	BC	162	GLN	3.3
29	BE	149	ILE	3.3
29	DE	73	ILE	3.3
37	DL	142	ILE	3.3
5	AF	55	HIS	3.3
11	AL	92	VAL	3.3
13	AP	5	ARG	3.3
25	DC	15	VAL	3.3
26	BD	26	VAL	3.3
26	DD	72	GLY	3.3
16	AS	47	THR	3.3
29	BE	15	SER	3.3
52	BW	63	ASP	3.3
26	BD	209	ALA	3.3
5	AF	84	VAL	3.3
27	BK	120	PRO	3.3
37	DL	126	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
24	BI	98	GLY	3.3
51	BZ	49	LEU	3.3
25	BC	101	ARG	3.3
27	DK	9	ASN	3.3
50	DT	3	ARG	3.3
24	DI	44	LYS	3.3
37	DL	107	PHE	3.3
45	DS	75	PHE	3.3
47	DF	37	MET	3.3
13	AP	60	TRP	3.3
27	BK	53	LYS	3.3
37	DL	96	LYS	3.3
38	BM	75	GLU	3.3
25	DC	91	ALA	3.3
41	BJ	66	GLY	3.3
47	BF	116	LEU	3.3
42	BN	10	LEU	3.3
47	DF	29	ARG	3.3
25	BC	171	VAL	3.3
37	DL	119	PRO	3.3
27	BK	75	SER	3.3
44	DQ	117	ALA	3.3
4	CE	110	MET	3.3
25	BC	100	ARG	3.3
25	BC	153	LEU	3.3
27	DK	84	CYS	3.3
43	DO	37	ALA	3.3
21	CN	46	LYS	3.3
47	DF	16	MET	3.3
5	AF	71	ILE	3.3
12	AM	52	ILE	3.3
38	DM	117	PHE	3.3
4	CE	155	LYS	3.3
4	AE	85	LYS	3.2
7	CH	116	ARG	3.2
20	AO	46	HIS	3.2
21	AN	27	LYS	3.2
41	BJ	72	LYS	3.2
11	CL	24	GLU	3.2
49	DR	50	GLY	3.2
24	BI	108	ILE	3.2
37	DL	110	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
50	DT	47	VAL	3.2
12	AM	92	ARG	3.2
25	BC	102	TYR	3.2
29	DE	57	LYS	3.2
24	BI	8	VAL	3.2
37	DL	46	VAL	3.2
9	AJ	62	ARG	3.2
25	DC	60	ALA	3.2
40	BH	89	LYS	3.2
24	BI	140	GLU	3.2
38	BM	67	VAL	3.2
21	AN	30	ILE	3.2
51	BZ	60	ASP	3.2
48	BG	93	TYR	3.2
18	CB	197	PHE	3.2
18	CB	30	ILE	3.2
40	DH	80	ILE	3.2
47	BF	35	LEU	3.2
45	DS	89	ALA	3.2
4	CE	86	GLY	3.2
23	BB	2309	A	3.2
2	CC	166	TRP	3.2
7	AH	98	LEU	3.2
24	DI	125	THR	3.2
3	AD	174	ALA	3.2
34	D3	20	GLY	3.2
46	BU	17	ASP	3.2
14	AQ	74	LEU	3.2
47	DF	2	LYS	3.2
24	DI	53	PRO	3.2
50	DT	55	VAL	3.2
27	DK	17	ARG	3.2
18	CB	38	HIS	3.2
13	AP	29	ASN	3.2
14	AQ	33	TYR	3.2
27	DK	45	GLU	3.2
43	DO	115	LEU	3.2
48	BG	132	LEU	3.2
24	BI	41	PHE	3.2
35	DV	72	VAL	3.2
40	BH	135	HIS	3.2
5	AF	12	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
5	AF	95	ALA	3.2
9	CJ	74	VAL	3.2
25	DC	19	VAL	3.2
29	DE	121	VAL	3.2
35	BV	61	LEU	3.2
52	DW	36	ILE	3.2
18	CB	29	PHE	3.2
19	AU	4	LYS	3.2
25	BC	124	LYS	3.2
27	BK	56	ASP	3.2
13	AP	6	LEU	3.2
16	AS	30	LEU	3.2
36	D2	31	LEU	3.2
5	AF	89	VAL	3.2
47	BF	174	PHE	3.2
23	DB	1731	G	3.1
25	BC	145	MET	3.1
7	CH	122	GLY	3.1
21	AN	65	GLN	3.1
26	DD	77	ARG	3.1
40	BH	97	ARG	3.1
6	AG	73	GLU	3.1
4	CE	38	VAL	3.1
5	AF	80	PHE	3.1
28	BP	62	LYS	3.1
38	DM	108	VAL	3.1
35	DV	94	ALA	3.1
9	CJ	33	GLY	3.1
37	DL	95	LEU	3.1
35	DV	56	PHE	3.1
46	BU	72	PHE	3.1
5	AF	61	LEU	3.1
27	BK	99	ILE	3.1
36	B2	31	LEU	3.1
38	BM	7	THR	3.1
14	AQ	77	VAL	3.1
4	CE	119	VAL	3.1
26	DD	34	VAL	3.1
41	BJ	75	TYR	3.1
9	CJ	80	THR	3.1
25	BC	50	THR	3.1
35	BV	55	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
29	BE	129	PRO	3.1
11	AL	16	ALA	3.1
28	BP	69	VAL	3.1
39	BX	24	GLU	3.1
21	AN	51	PRO	3.1
25	DC	21	PRO	3.1
5	AF	56	LYS	3.1
9	CJ	30	LYS	3.1
25	DC	204	LEU	3.1
47	DF	14	LYS	3.1
18	CB	186	VAL	3.1
24	BI	23	VAL	3.1
18	CB	70	GLY	3.1
46	BU	86	PHE	3.1
39	DX	13	GLU	3.1
35	BV	83	LYS	3.1
2	AC	67	ILE	3.1
29	DE	125	SER	3.1
32	B4	23	ILE	3.1
5	AF	21	MET	3.1
47	DF	166	ARG	3.1
27	BK	82	ASN	3.1
25	DC	56	GLY	3.1
27	DK	83	ALA	3.1
13	AP	9	HIS	3.1
16	AS	74	ALA	3.1
27	BK	83	ALA	3.1
40	BH	108	VAL	3.1
28	DP	1	SER	3.1
37	DL	109	LYS	3.1
9	CJ	91	ASP	3.1
23	BB	1095	A	3.1
28	BP	40	GLN	3.1
12	CM	7	ASN	3.1
37	DL	60	ARG	3.1
18	CB	67	LEU	3.1
38	DM	133	LYS	3.1
9	CJ	76	ILE	3.1
25	BC	5	CYS	3.1
15	CR	31	TYR	3.1
47	BF	176	PHE	3.1
15	AR	66	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
26	BD	110	THR	3.1
26	DD	4	LEU	3.1
43	DO	23	ALA	3.0
40	BH	114	GLU	3.0
27	BK	95	ILE	3.0
45	DS	74	ILE	3.0
11	AL	15	VAL	3.0
21	CN	21	ALA	3.0
24	DI	43	ALA	3.0
34	D3	4	LYS	3.0
37	DL	75	ALA	3.0
24	DI	49	GLU	3.0
48	BG	37	ASN	3.0
14	AQ	9	GLY	3.0
26	BD	140	HIS	3.0
5	AF	81	ASN	3.0
20	AO	3	LEU	3.0
16	AS	75	PRO	3.0
35	DV	7	GLU	3.0
39	DX	1	MET	3.0
27	BK	85	VAL	3.0
12	AM	51	GLN	3.0
20	AO	47	LYS	3.0
44	DQ	90	ASP	3.0
7	CH	73	SER	3.0
25	BC	92	LEU	3.0
18	CB	99	MET	3.0
2	CC	203	LYS	3.0
4	CE	71	ILE	3.0
6	AG	13	PRO	3.0
47	DF	84	ILE	3.0
25	BC	167	ASP	3.0
12	AM	83	GLY	3.0
27	BK	105	ARG	3.0
36	D2	41	ARG	3.0
45	DS	95	ARG	3.0
38	DM	67	VAL	3.0
29	BE	154	ASP	3.0
38	DM	102	LEU	3.0
43	DO	62	LEU	3.0
9	CJ	27	GLU	3.0
27	BK	63	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
27	BK	106	GLU	3.0
45	DS	110	ARG	3.0
40	DH	99	ILE	3.0
49	BR	48	LYS	3.0
23	BB	62	U	3.0
27	BK	86	LEU	3.0
47	BF	151	LEU	3.0
39	DX	23	ARG	3.0
47	DF	9	ASP	3.0
3	CD	106	PHE	3.0
2	CC	181	ILE	3.0
21	AN	56	PRO	3.0
21	CN	24	ALA	3.0
24	BI	86	LYS	3.0
25	BC	2	VAL	3.0
38	BM	90	GLU	3.0
26	DD	76	GLY	3.0
38	BM	31	PHE	3.0
34	D3	53	ASP	3.0
46	BU	87	GLU	3.0
47	DF	3	LEU	3.0
4	CE	111	ARG	3.0
9	AJ	34	ALA	3.0
16	AS	10	ILE	3.0
5	AF	58	HIS	3.0
40	BH	5	LEU	3.0
50	DT	32	LEU	3.0
4	CE	118	GLY	3.0
11	AL	12	ALA	3.0
29	BE	11	ALA	3.0
29	DE	54	GLY	3.0
50	DT	80	TRP	3.0
1	AA	121	U	3.0
7	CH	102	VAL	3.0
16	AS	48	ILE	3.0
48	BG	91	VAL	3.0
29	DE	106	LYS	3.0
37	DL	125	LEU	3.0
40	BH	27	ARG	3.0
18	CB	52	ALA	3.0
1	AA	1229	A	3.0
4	CE	84	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
23	DB	752	A	3.0
26	DD	56	LYS	3.0
13	AP	17	TYR	3.0
41	BJ	74	TYR	3.0
12	AM	59	VAL	2.9
2	AC	100	ILE	2.9
25	DC	222	THR	2.9
3	CD	202	LEU	2.9
29	DE	49	ARG	2.9
24	DI	82	ALA	2.9
25	BC	95	TYR	2.9
29	DE	108	ILE	2.9
47	BF	103	ILE	2.9
40	BH	139	PHE	2.9
3	AD	175	GLY	2.9
11	AL	91	GLY	2.9
13	AP	19	VAL	2.9
27	DK	50	GLY	2.9
43	BO	89	ASP	2.9
39	BX	1	MET	2.9
47	DF	94	ARG	2.9
50	DT	61	LEU	2.9
14	AQ	27	PHE	2.9
49	BR	35	PHE	2.9
48	BG	21	GLN	2.9
48	DG	59	ASP	2.9
50	DT	52	GLU	2.9
16	CS	25	GLY	2.9
47	DF	159	ALA	2.9
6	AG	46	LEU	2.9
21	AN	58	ARG	2.9
48	BG	59	ASP	2.9
40	BH	96	THR	2.9
47	DF	168	LEU	2.9
16	AS	68	HIS	2.9
21	CN	76	PHE	2.9
35	DV	47	VAL	2.9
45	DS	83	LYS	2.9
18	AB	100	LEU	2.9
6	AG	150	PHE	2.9
9	CJ	88	MET	2.9
11	AL	24	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
21	AN	91	GLU	2.9
23	BB	1733	G	2.9
38	BM	3	GLN	2.9
25	DC	215	VAL	2.9
1	AA	1230	C	2.9
28	BP	63	ILE	2.9
42	BN	48	VAL	2.9
2	CC	186	SER	2.9
25	BC	109	LEU	2.9
2	CC	150	VAL	2.9
50	DT	54	GLU	2.9
21	AN	47	LEU	2.9
25	BC	104	LEU	2.9
37	DL	93	ASN	2.9
47	DF	45	ASP	2.9
35	BV	2	PHE	2.9
40	DH	110	VAL	2.9
29	DE	92	HIS	2.9
42	BN	17	ARG	2.9
9	AJ	63	ASP	2.9
25	DC	18	VAL	2.9
34	D3	22	LYS	2.9
40	BH	48	GLU	2.9
21	AN	48	GLN	2.9
24	BI	5	GLN	2.9
24	BI	79	LEU	2.9
25	BC	103	ILE	2.9
2	AC	151	GLU	2.9
25	BC	179	GLU	2.9
27	BK	7	MET	2.9
40	DH	17	ASP	2.9
27	BK	108	ARG	2.9
7	AH	60	LEU	2.9
14	AQ	52	CYS	2.9
45	DS	105	VAL	2.9
3	CD	160	LEU	2.9
8	AI	83	THR	2.9
25	BC	163	ILE	2.9
26	BD	27	ILE	2.9
25	BC	72	GLY	2.9
47	DF	165	GLY	2.9
16	AS	58	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
50	DT	16	VAL	2.9
34	D3	12	ARG	2.9
21	CN	29	ILE	2.9
28	BP	65	ASN	2.9
4	AE	95	MET	2.8
50	DT	4	GLU	2.8
2	AC	156	LEU	2.8
6	AG	14	ASP	2.8
12	AM	38	ILE	2.8
18	CB	53	LEU	2.8
38	BM	2	LEU	2.8
26	BD	109	VAL	2.8
34	D3	57	VAL	2.8
34	D3	60	CYS	2.8
37	DL	48	ARG	2.8
4	CE	75	LEU	2.8
21	AN	26	LEU	2.8
36	D2	25	LYS	2.8
36	D2	45	SER	2.8
40	DH	72	ILE	2.8
24	BI	44	LYS	2.8
25	BC	70	LYS	2.8
27	DK	82	ASN	2.8
42	BN	86	ARG	2.8
27	DK	43	ILE	2.8
37	DL	122	VAL	2.8
40	DH	140	ALA	2.8
25	DC	20	ASN	2.8
25	DC	198	GLU	2.8
7	CH	44	PHE	2.8
12	AM	32	ILE	2.8
28	BP	98	TYR	2.8
40	BH	143	ILE	2.8
26	DD	94	GLN	2.8
28	BP	27	VAL	2.8
39	BX	10	SER	2.8
47	BF	131	VAL	2.8
2	CC	139	ASN	2.8
18	AB	66	ILE	2.8
24	BI	10	LEU	2.8
29	DE	172	ALA	2.8
41	DJ	20	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
27	BK	3	GLN	2.8
47	DF	176	PHE	2.8
18	CB	157	PRO	2.8
27	DK	42	THR	2.8
28	BP	45	VAL	2.8
34	B3	19	GLY	2.8
4	CE	143	LEU	2.8
7	CH	91	LEU	2.8
21	AN	76	PHE	2.8
27	BK	107	LEU	2.8
47	DF	25	MET	2.8
51	BZ	38	PHE	2.8
35	DV	57	TYR	2.8
48	BG	167	VAL	2.8
41	DJ	63	ALA	2.8
16	CS	19	GLU	2.8
21	AN	31	SER	2.8
23	BB	613	A	2.8
23	DB	846	U	2.8
40	BH	107	GLY	2.8
28	BP	86	LYS	2.8
29	DE	48	THR	2.8
16	AS	64	GLU	2.8
23	DB	1728	C	2.8
27	BK	39	ILE	2.8
38	DM	126	ILE	2.8
8	AI	48	ARG	2.8
25	BC	143	VAL	2.8
36	D2	28	ARG	2.8
2	AC	44	LYS	2.8
48	DG	90	GLY	2.8
16	CS	47	THR	2.8
29	DE	51	GLU	2.8
25	BC	110	LYS	2.8
29	BE	172	ALA	2.8
50	DT	83	ALA	2.8
29	BE	175	ILE	2.8
12	AM	78	ARG	2.8
40	DH	9	VAL	2.8
43	DO	39	VAL	2.8
29	DE	56	GLY	2.8
7	AH	44	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
16	AS	70	LEU	2.8
47	BF	56	LEU	2.8
8	AI	20	ILE	2.8
38	DM	35	ALA	2.8
9	CJ	90	LEU	2.8
14	AQ	43	LEU	2.8
48	DG	147	LEU	2.8
3	AD	177	MET	2.8
28	BP	25	VAL	2.8
11	AL	87	LYS	2.8
41	BJ	65	THR	2.8
24	BI	66	PHE	2.8
47	BF	120	SER	2.8
13	AP	34	GLU	2.8
47	BF	71	LYS	2.8
2	AC	46	LEU	2.8
12	AM	104	ASN	2.8
28	BP	59	THR	2.8
45	DS	97	LEU	2.8
47	DF	15	LEU	2.8
2	CC	93	ILE	2.7
21	AN	60	ARG	2.7
37	DL	90	VAL	2.7
13	AP	38	PHE	2.7
18	CB	49	PHE	2.7
23	BB	136	G	2.7
18	CB	124	THR	2.7
38	DM	107	GLY	2.7
44	DQ	17	LEU	2.7
41	BJ	68	LYS	2.7
47	BF	142	TYR	2.7
25	DC	101	ARG	2.7
29	DE	150	THR	2.7
9	CJ	29	ALA	2.7
21	CN	45	LEU	2.7
2	CC	130	ARG	2.7
25	DC	65	ASP	2.7
40	BH	61	VAL	2.7
18	CB	213	LEU	2.7
38	DM	40	ARG	2.7
25	DC	229	HIS	2.7
36	D2	42	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
48	BG	51	PHE	2.7
25	BC	75	ALA	2.7
37	DL	83	ALA	2.7
3	CD	34	GLU	2.7
21	CN	25	GLU	2.7
25	BC	78	GLU	2.7
21	AN	34	ASN	2.7
26	DD	180	VAL	2.7
41	BJ	84	ILE	2.7
25	BC	173	LEU	2.7
25	DC	61	TYR	2.7
4	CE	44	ARG	2.7
9	CJ	6	ILE	2.7
47	DF	78	ILE	2.7
25	BC	99	GLU	2.7
35	DV	42	LEU	2.7
38	DM	59	ARG	2.7
4	CE	82	HIS	2.7
5	AF	92	THR	2.7
48	BG	122	ALA	2.7
26	BD	148	GLN	2.7
48	BG	87	GLN	2.7
49	DR	48	LYS	2.7
52	DW	43	LYS	2.7
13	AP	37	GLY	2.7
25	BC	137	GLY	2.7
45	BS	21	ALA	2.7
7	CH	62	LEU	2.7
18	CB	57	ASN	2.7
25	BC	94	LEU	2.7
38	DM	136	MET	2.7
7	AH	128	VAL	2.7
24	BI	31	GLY	2.7
25	DC	90	ILE	2.7
21	CN	47	LEU	2.7
26	BD	28	GLU	2.7
26	DD	74	GLU	2.7
4	CE	45	VAL	2.7
25	BC	183	VAL	2.7
27	BK	101	GLY	2.7
48	BG	160	GLY	2.7
25	BC	138	SER	2.7

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Mol	Chain	Res	Type	RSRZ
26	DD	30	GLU	2.7
29	DE	30	GLN	2.7
39	DX	24	GLU	2.7
51	BZ	71	LEU	2.7
33	D1	4	ILE	2.7
40	BH	106	ALA	2.7
2	CC	110	LEU	2.7
2	CC	182	ASP	2.7
3	CD	169	TRP	2.7
12	AM	86	ARG	2.7
39	BX	9	LYS	2.7
47	DF	56	LEU	2.7
47	DF	162	ASP	2.7
26	DD	203	VAL	2.7
27	DK	10	VAL	2.7
28	BP	74	GLN	2.7
29	DE	126	VAL	2.7
16	AS	9	PHE	2.7
19	AU	36	PHE	2.7
2	CC	151	GLU	2.7
26	BD	181	ASP	2.7
52	BW	82	GLU	2.7
14	AQ	6	THR	2.7
15	AR	63	TYR	2.7
23	DB	2799	A	2.7
3	CD	181	PHE	2.7
16	CS	40	PHE	2.7
25	DC	36	ASN	2.7
27	DK	51	LYS	2.7
35	DV	73	LYS	2.7
50	DT	49	LYS	2.7
25	DC	92	LEU	2.6
7	CH	95	MET	2.6
25	BC	64	VAL	2.6
26	DD	17	GLU	2.6
37	DL	44	GLY	2.6
40	BH	35	LYS	2.6
48	BG	114	HIS	2.6
23	DB	1729	U	2.6
25	BC	132	ARG	2.6
28	BP	91	VAL	2.6
37	DL	100	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
5	AF	32	ALA	2.6
29	DE	190	ALA	2.6
9	CJ	73	LEU	2.6
40	BH	77	THR	2.6
12	AM	7	ASN	2.6
16	AS	72	GLU	2.6
39	BX	60	LYS	2.6
16	AS	26	ASP	2.6
19	AU	26	GLY	2.6
29	DE	90	GLN	2.6
47	DF	117	SER	2.6
16	AS	78	THR	2.6
29	BE	169	VAL	2.6
8	AI	29	ILE	2.6
25	DC	202	ARG	2.6
47	DF	151	LEU	2.6
7	AH	102	VAL	2.6
29	BE	121	VAL	2.6
37	DL	56	PRO	2.6
45	DS	4	ILE	2.6
27	BK	9	ASN	2.6
14	AQ	22	VAL	2.6
24	DI	115	ASP	2.6
34	D3	49	VAL	2.6
40	BH	136	SER	2.6
21	AN	46	LYS	2.6
24	DI	70	THR	2.6
50	BT	5	GLU	2.6
6	AG	78	ARG	2.6
48	BG	92	GLY	2.6
21	CN	54	SER	2.6
14	AQ	56	ASP	2.6
5	AF	59	TYR	2.6
6	AG	57	GLU	2.6
36	D2	39	ARG	2.6
16	CS	29	PRO	2.6
50	BT	93	LEU	2.6
25	DC	64	VAL	2.6
32	B4	25	VAL	2.6
4	CE	42	ASN	2.6
7	CH	74	ILE	2.6
18	CB	39	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
18	CB	90	PHE	2.6
25	DC	212	TRP	2.6
18	CB	159	ALA	2.6
24	DI	131	THR	2.6
29	DE	105	LEU	2.6
25	DC	17	LYS	2.6
16	AS	31	ARG	2.6
25	BC	166	ARG	2.6
29	DE	42	GLY	2.6
40	DH	5	LEU	2.6
41	BJ	122	LEU	2.6
34	D3	64	ALA	2.6
27	DK	3	GLN	2.6
34	D3	27	ASN	2.6
4	CE	55	VAL	2.6
41	BJ	105	VAL	2.6
15	AR	60	ARG	2.6
37	DL	78	ARG	2.6
9	AJ	65	TYR	2.6
42	BN	87	PHE	2.6
16	CS	65	MET	2.6
25	DC	80	LEU	2.6
21	CN	53	ASP	2.6
29	DE	168	ASP	2.6
48	DG	53	PRO	2.6
42	BN	82	GLU	2.6
50	DT	74	ILE	2.6
33	B1	49	LYS	2.6
45	DS	82	MET	2.6
45	DS	90	LYS	2.6
18	CB	75	ALA	2.6
20	AO	44	ALA	2.6
24	DI	114	ALA	2.6
48	BG	90	GLY	2.6
40	DH	38	PRO	2.6
13	CP	45	GLU	2.6
29	BE	152	GLU	2.6
48	DG	106	LEU	2.6
4	CE	122	VAL	2.6
11	AL	60	PHE	2.6
18	CB	34	ARG	2.6
11	AL	62	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
14	AQ	57	VAL	2.6
29	BE	187	VAL	2.6
37	DL	116	VAL	2.6
48	DG	56	GLY	2.6
29	DE	93	SER	2.6
35	DV	93	ARG	2.6
21	AN	25	GLU	2.5
29	DE	158	PHE	2.5
38	BM	77	PRO	2.5
47	BF	64	PRO	2.5
18	CB	45	THR	2.5
18	AB	161	PHE	2.5
47	DF	99	PHE	2.5
10	AK	128	VAL	2.5
1	AA	461	A	2.5
25	BC	125	PRO	2.5
2	AC	99	GLN	2.5
21	AN	52	ARG	2.5
36	B2	21	ARG	2.5
45	BS	94	ASP	2.5
9	CJ	83	THR	2.5
50	DT	45	ALA	2.5
2	CC	200	TRP	2.5
7	AH	101	ALA	2.5
25	BC	121	ALA	2.5
38	DM	62	LYS	2.5
1	AA	1228	C	2.5
35	DV	4	ILE	2.5
48	BG	102	ILE	2.5
14	AQ	11	VAL	2.5
25	DC	2	VAL	2.5
41	BJ	62	VAL	2.5
5	AF	79	ARG	2.5
27	BK	40	LYS	2.5
27	BK	54	LYS	2.5
52	DW	65	LYS	2.5
27	DK	8	LEU	2.5
35	DV	33	GLY	2.5
23	DB	645	C	2.5
2	CC	69	THR	2.5
25	BC	86	ARG	2.5
4	CE	149	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
5	AF	50	PRO	2.5
11	AL	18	SER	2.5
29	BE	151	GLY	2.5
41	DJ	4	PHE	2.5
47	DF	77	LYS	2.5
47	BF	135	ILE	2.5
50	DT	8	LEU	2.5
5	AF	82	ASP	2.5
11	AL	17	LYS	2.5
25	BC	172	THR	2.5
45	DS	86	MET	2.5
13	AP	54	LEU	2.5
24	BI	57	VAL	2.5
25	BC	21	PRO	2.5
35	DV	63	ILE	2.5
35	BV	93	ARG	2.5
42	BN	20	MET	2.5
5	AF	57	ALA	2.5
5	AF	60	VAL	2.5
21	AN	54	SER	2.5
26	DD	146	ILE	2.5
27	BK	100	PHE	2.5
37	DL	84	LYS	2.5
47	DF	6	TYR	2.5
37	DL	111	ILE	2.5
38	DM	57	VAL	2.5
47	DF	96	TRP	2.5
39	DX	17	GLU	2.5
41	BJ	20	ALA	2.5
48	BG	116	LEU	2.5
3	AD	145	ARG	2.5
25	BC	142	ASN	2.5
47	DF	161	SER	2.5
27	BK	112	PHE	2.5
47	DF	31	GLU	2.5
11	AL	8	ARG	2.5
18	CB	33	ALA	2.5
38	DM	135	VAL	2.5
21	AN	59	GLN	2.5
38	BM	104	GLU	2.5
25	DC	214	GLY	2.5
26	DD	84	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
48	BG	20	GLY	2.5
9	CJ	75	ASP	2.5
25	BC	106	PRO	2.5
26	DD	101	PHE	2.5
27	BK	79	PHE	2.5
47	BF	152	ASP	2.5
29	DE	62	GLN	2.5
37	DL	106	GLU	2.5
18	AB	183	PHE	2.5
34	D3	63	TYR	2.4
11	AL	56	LEU	2.4
37	DL	74	THR	2.4
37	DL	115	GLU	2.4
38	BM	128	THR	2.4
47	DF	35	LEU	2.4
5	AF	85	ILE	2.4
7	CH	100	ILE	2.4
25	BC	165	ALA	2.4
26	BD	192	ALA	2.4
27	BK	60	ALA	2.4
38	BM	40	ARG	2.4
41	BJ	101	ILE	2.4
47	BF	111	ARG	2.4
4	CE	47	PHE	2.4
7	CH	112	ASP	2.4
26	BD	161	MET	2.4
41	DJ	24	THR	2.4
44	DQ	36	GLN	2.4
8	AI	126	PHE	2.4
25	BC	117	SER	2.4
25	BC	126	GLY	2.4
27	BK	81	GLY	2.4
36	D2	4	THR	2.4
3	CD	89	LEU	2.4
20	AO	67	LEU	2.4
28	BP	102	ARG	2.4
43	DO	48	LEU	2.4
5	AF	96	VAL	2.4
46	BU	91	LYS	2.4
27	DK	2	ILE	2.4
23	BB	1459	G	2.4
4	CE	28	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
11	AL	51	VAL	2.4
47	DF	163	GLU	2.4
46	BU	102	ILE	2.4
32	B4	35	GLN	2.4
27	BK	109	SER	2.4
28	BP	64	SER	2.4
45	BS	104	THR	2.4
4	CE	76	ASN	2.4
14	AQ	7	LEU	2.4
19	AU	34	ARG	2.4
27	DK	63	VAL	2.4
38	DM	110	GLU	2.4
50	DT	42	GLU	2.4
50	DT	64	LYS	2.4
23	DB	204	A	2.4
47	BF	43	ILE	2.4
8	AI	125	GLN	2.4
18	AB	158	ASP	2.4
21	AN	95	LEU	2.4
43	DO	69	ASP	2.4
25	BC	76	VAL	2.4
25	DC	93	VAL	2.4
40	DH	147	VAL	2.4
28	BP	76	HIS	2.4
29	BE	19	PHE	2.4
42	BN	25	ALA	2.4
47	DF	178	LYS	2.4
3	CD	162	GLU	2.4
40	DH	76	GLU	2.4
6	AG	6	ILE	2.4
7	AH	100	ILE	2.4
27	BK	34	GLY	2.4
29	DE	119	ILE	2.4
23	DB	1459	G	2.4
18	AB	42	LEU	2.4
4	CE	32	PHE	2.4
14	AQ	42	LYS	2.4
6	AG	4	ARG	2.4
23	DB	789	A	2.4
24	DI	80	LYS	2.4
29	BE	6	LYS	2.4
25	BC	155	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
14	CQ	12	VAL	2.4
28	BP	60	VAL	2.4
40	BH	18	GLN	2.4
27	BK	23	LYS	2.4
35	DV	3	THR	2.4
25	BC	91	ALA	2.4
47	BF	98	PHE	2.4
3	CD	107	GLY	2.4
21	AN	57	SER	2.4
44	DQ	116	LEU	2.4
51	BZ	40	VAL	2.4
25	BC	118	GLY	2.4
25	DC	205	GLY	2.4
2	CC	111	ASP	2.4
12	CM	47	LEU	2.4
27	DK	64	ARG	2.4
45	DS	108	SER	2.4
32	B4	22	VAL	2.4
48	DG	42	VAL	2.4
25	BC	74	PRO	2.4
21	CN	27	LYS	2.4
26	BD	15	PHE	2.4
43	DO	61	GLN	2.4
13	AP	51	ARG	2.4
18	CB	16	GLY	2.4
23	DB	791	C	2.4
45	DS	9	HIS	2.4
16	CS	30	LEU	2.4
20	AO	66	LEU	2.4
29	BE	12	LEU	2.4
16	AS	63	ASP	2.4
40	DH	19	VAL	2.4
46	BU	82	VAL	2.4
2	AC	187	GLU	2.4
3	CD	190	LEU	2.4
36	B2	32	ALA	2.4
39	BX	57	LEU	2.4
45	DS	51	LEU	2.4
12	AM	103	THR	2.4
28	DP	37	LYS	2.4
34	D3	3	ILE	2.3
34	D3	14	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
48	BG	38	ASP	2.4
16	CS	48	ILE	2.3
37	DL	135	ILE	2.3
40	DH	94	ILE	2.3
29	DE	78	TRP	2.3
48	BG	169	ARG	2.3
3	AD	24	VAL	2.3
7	AH	58	LEU	2.3
11	AL	52	CYS	2.3
38	DM	41	LEU	2.3
49	DR	75	VAL	2.3
13	AP	4	ILE	2.3
27	BK	51	LYS	2.3
32	B4	8	LYS	2.3
43	DO	40	ILE	2.3
24	DI	135	MET	2.3
25	DC	219	VAL	2.3
41	BJ	59	ALA	2.3
41	BJ	67	ASN	2.3
47	DF	69	ALA	2.3
2	CC	146	LYS	2.3
16	CS	24	SER	2.3
13	AP	55	ASP	2.3
27	BK	78	ARG	2.3
6	AG	70	PRO	2.3
10	AK	73	VAL	2.3
13	AP	71	VAL	2.3
25	BC	168	GLY	2.3
43	BO	114	GLY	2.3
27	BK	11	ALA	2.3
25	DC	100	ARG	2.3
24	DI	88	GLY	2.3
40	DH	75	LEU	2.3
48	DG	113	ASP	2.3
7	CH	101	ALA	2.3
25	BC	29	PHE	2.3
2	CC	201	ILE	2.3
25	DC	42	ARG	2.3
27	BK	43	ILE	2.3
27	DK	38	ILE	2.3
45	DS	96	ILE	2.3
51	BZ	64	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
28	BP	26	GLU	2.3
35	BV	59	GLU	2.3
26	BD	130	GLN	2.3
13	AP	39	PHE	2.3
42	DN	21	PHE	2.3
23	BB	545	U	2.3
7	AH	59	GLU	2.3
27	BK	61	VAL	2.3
43	BO	26	LEU	2.3
3	CD	144	ILE	2.3
17	CT	38	ILE	2.3
27	DK	39	ILE	2.3
41	BJ	69	ARG	2.3
18	CB	43	GLU	2.3
40	DH	71	LYS	2.3
50	DT	58	VAL	2.3
18	CB	15	PHE	2.3
21	CN	16	ALA	2.3
23	BB	878	A	2.3
28	DP	19	PHE	2.3
47	BF	74	ALA	2.3
7	CH	124	ILE	2.3
25	BC	90	ILE	2.3
38	DM	114	ARG	2.3
25	BC	3	VAL	2.3
28	DP	67	GLU	2.3
5	AF	38	ARG	2.3
11	AL	25	ALA	2.3
24	DI	20	SER	2.3
36	D2	21	ARG	2.3
38	DM	63	ILE	2.3
16	AS	50	VAL	2.3
26	BD	189	VAL	2.3
28	BP	96	LEU	2.3
40	BH	9	VAL	2.3
40	BH	83	LYS	2.3
43	DO	53	THR	2.3
48	DG	116	LEU	2.3
48	DG	161	VAL	2.3
2	CC	94	ALA	2.3
32	B4	24	ARG	2.3
48	DG	54	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
15	AR	25	ILE	2.3
21	AN	93	PRO	2.3
41	BJ	55	ILE	2.3
21	CN	97	LYS	2.3
9	CJ	28	THR	2.3
24	BI	67	THR	2.3
40	BH	137	GLU	2.3
25	DC	231	HIS	2.3
44	DQ	32	ARG	2.3
48	DG	58	ALA	2.3
18	CB	40	ILE	2.3
29	DE	181	ILE	2.3
29	DE	59	PRO	2.3
43	BO	78	VAL	2.3
25	BC	184	GLU	2.3
35	DV	69	GLU	2.3
4	CE	15	ILE	2.3
2	AC	90	VAL	2.3
34	D3	28	LEU	2.3
12	CM	9	PRO	2.3
48	BG	172	GLU	2.3
19	AU	16	ARG	2.3
28	BP	68	GLY	2.3
36	B2	35	ARG	2.3
47	BF	34	THR	2.3
25	BC	264	LYS	2.3
29	BE	57	LYS	2.3
2	CC	109	GLU	2.3
25	BC	66	PHE	2.3
30	BY	58	GLU	2.3
38	BM	4	PRO	2.3
45	DS	80	PRO	2.3
47	DF	150	GLY	2.3
5	AF	36	ILE	2.3
4	CE	113	VAL	2.3
7	CH	128	VAL	2.3
51	BZ	47	VAL	2.3
20	AO	4	SER	2.3
25	DC	29	PHE	2.3
47	DF	20	ASN	2.3
13	AP	33	ILE	2.3
14	AQ	78	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
16	AS	14	LEU	2.3
20	AO	75	VAL	2.3
27	BK	62	VAL	2.3
41	BJ	17	VAL	2.3
5	AF	75	GLU	2.2
7	CH	126	CYS	2.2
18	AB	192	PRO	2.2
20	AO	6	GLU	2.2
23	DB	2609	U	2.2
27	DK	21	CYS	2.2
28	DP	40	GLN	2.2
30	DY	58	GLU	2.2
28	BP	44	GLY	2.2
25	BC	85	ASN	2.2
27	DK	65	THR	2.2
25	BC	141	HIS	2.2
2	AC	63	ILE	2.2
24	DI	35	MET	2.2
29	DE	178	VAL	2.2
10	AK	80	ASN	2.2
5	AF	86	ARG	2.2
10	AK	126	ARG	2.2
24	BI	125	THR	2.2
40	BH	17	ASP	2.2
42	BN	78	LYS	2.2
49	BR	45	GLU	2.2
2	CC	70	ALA	2.2
4	CE	97	PRO	2.2
37	DL	124	GLY	2.2
5	AF	29	ILE	2.2
16	AS	46	LEU	2.2
45	DS	21	ALA	2.2
47	BF	140	ILE	2.2
8	AI	40	ARG	2.2
25	BC	62	ARG	2.2
37	DL	2	ARG	2.2
31	B0	56	LYS	2.2
13	AP	18	GLN	2.2
45	DS	109	ASP	2.2
8	CI	51	LEU	2.2
40	BH	15	LEU	2.2
41	BJ	73	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
48	BG	53	PRO	2.2
7	AH	127	TYR	2.2
29	DE	23	PHE	2.2
29	DE	44	ARG	2.2
4	CE	9	GLU	2.2
7	CH	94	VAL	2.2
8	AI	55	ASP	2.2
18	CB	48	MET	2.2
19	AU	27	VAL	2.2
23	BB	764	A	2.2
23	DB	2602	A	2.2
24	DI	79	LEU	2.2
28	BP	72	VAL	2.2
45	BS	85	ILE	2.2
45	DS	107	VAL	2.2
5	AF	49	TYR	2.2
23	DB	1727	C	2.2
23	DB	1781	U	2.2
26	BD	118	PHE	2.2
36	D2	34	ARG	2.2
44	DQ	5	ARG	2.2
42	BN	56	LYS	2.2
29	DE	81	GLY	2.2
28	BP	106	ALA	2.2
40	DH	4	ILE	2.2
25	BC	174	ARG	2.2
26	BD	156	PHE	2.2
1	AA	86	G	2.2
15	CR	32	ILE	2.2
2	AC	98	ALA	2.2
16	AS	80	ARG	2.2
29	BE	61	ARG	2.2
30	DY	2	LYS	2.2
3	CD	60	VAL	2.2
7	CH	120	LEU	2.2
8	AI	54	VAL	2.2
27	BK	115	ILE	2.2
47	BF	153	ILE	2.2
50	DT	85	VAL	2.2
3	AD	104	MET	2.2
5	AF	44	ARG	2.2
14	AQ	80	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
26	DD	31	ALA	2.2
43	BO	117	PHE	2.2
11	AL	10	PRO	2.2
22	BA	2	G	2.2
52	DW	45	HIS	2.2
3	AD	196	GLU	2.2
9	CJ	25	ILE	2.2
13	AP	75	ILE	2.2
43	BO	115	LEU	2.2
43	DO	28	VAL	2.2
8	AI	75	ALA	2.2
24	BI	109	ALA	2.2
25	BC	202	ARG	2.2
26	BD	132	ALA	2.2
47	BF	129	MET	2.2
47	DF	74	ALA	2.2
38	DM	3	GLN	2.2
6	AG	83	THR	2.2
27	BK	97	THR	2.2
2	CC	152	VAL	2.2
4	CE	142	GLY	2.2
26	BD	3	GLY	2.2
33	B1	34	GLU	2.2
27	BK	59	LYS	2.2
37	DL	52	GLY	2.2
40	BH	110	VAL	2.2
25	DC	250	GLN	2.2
35	BV	51	GLN	2.2
13	AP	53	ASP	2.2
26	DD	51	THR	2.2
1	CA	1030	U	2.2
25	BC	18	VAL	2.2
25	DC	16	VAL	2.2
25	DC	109	LEU	2.2
30	DY	56	VAL	2.2
40	BH	58	LEU	2.2
16	AS	17	LYS	2.2
23	DB	1566	A	2.2
27	BK	2	ILE	2.2
28	BP	49	ILE	2.2
27	BK	12	ASP	2.2
27	DK	12	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
38	DM	128	THR	2.2
51	BZ	50	ARG	2.2
45	DS	62	ASP	2.2
46	BU	90	LYS	2.2
41	DJ	6	ALA	2.2
25	DC	51	ARG	2.2
3	AD	107	GLY	2.2
28	BP	33	GLU	2.2
35	DV	84	PRO	2.2
39	BX	17	GLU	2.2
40	BH	49	ALA	2.2
20	AO	80	GLN	2.2
25	BC	119	VAL	2.2
25	DC	3	VAL	2.2
24	DI	48	ILE	2.2
26	DD	46	ARG	2.2
27	BK	24	VAL	2.2
27	BK	69	VAL	2.2
43	BO	88	LYS	2.2
43	DO	78	VAL	2.2
47	DF	26	GLN	2.2
40	DH	100	ALA	2.2
47	DF	52	ALA	2.2
28	BP	15	ASP	2.2
13	CP	54	LEU	2.2
13	CP	37	GLY	2.1
24	DI	140	GLU	2.2
24	BI	73	PRO	2.1
40	BH	13	GLY	2.1
8	AI	97	LEU	2.1
26	BD	201	LEU	2.1
42	BN	117	ASP	2.1
47	DF	79	ARG	2.1
34	D3	58	ILE	2.1
44	DQ	97	ILE	2.1
26	DD	57	ALA	2.1
27	BK	18	ARG	2.1
46	BU	92	VAL	2.1
24	BI	121	ILE	2.1
28	DP	75	THR	2.1
45	DS	24	ILE	2.1
14	CQ	8	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
9	CJ	39	PRO	2.1
27	DK	20	MET	2.1
16	AS	61	VAL	2.1
48	BG	168	VAL	2.1
2	CC	180	ASP	2.1
23	DB	1	G	2.1
16	CS	21	ALA	2.1
4	AE	122	VAL	2.1
5	AF	45	ARG	2.1
7	CH	103	VAL	2.1
14	AQ	10	ARG	2.1
37	DL	17	LYS	2.1
37	DL	89	VAL	2.1
28	BP	23	ASP	2.1
45	DS	7	HIS	2.1
47	BF	60	SER	2.1
18	CB	42	LEU	2.1
26	DD	5	VAL	2.1
41	BJ	56	VAL	2.1
48	DG	51	PHE	2.1
18	CB	162	VAL	2.1
26	DD	49	GLN	2.1
32	B4	9	LYS	2.1
27	DK	61	VAL	2.1
42	BN	12	ARG	2.1
11	AL	81	ILE	2.1
23	DB	613	A	2.1
24	BI	7	TYR	2.1
5	AF	33	GLU	2.1
47	DF	133	GLU	2.1
3	CD	177	MET	2.1
6	AG	86	VAL	2.1
8	AI	46	VAL	2.1
43	DO	26	LEU	2.1
47	DF	148	VAL	2.1
48	DG	89	VAL	2.1
25	BC	266	ILE	2.1
25	DC	28	PRO	2.1
28	BP	21	PRO	2.1
28	BP	83	ILE	2.1
36	D2	5	PHE	2.1
16	AS	45	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
40	DH	95	GLY	2.1
7	CH	60	LEU	2.1
29	BE	126	VAL	2.1
44	DQ	30	VAL	2.1
6	AG	82	SER	2.1
11	AL	39	THR	2.1
18	CB	14	HIS	2.1
47	BF	134	GLN	2.1
35	DV	82	TYR	2.1
8	AI	58	GLU	2.1
12	AM	87	GLY	2.1
25	BC	211	ARG	2.1
29	DE	180	LEU	2.1
42	DN	20	MET	2.1
2	AC	167	TYR	2.1
18	CB	188	THR	2.1
26	DD	97	SER	2.1
28	BP	84	SER	2.1
36	D2	24	THR	2.1
1	CA	1257	A	2.1
18	CB	41	ASN	2.1
2	AC	65	VAL	2.1
2	CC	169	GLU	2.1
27	DK	81	GLY	2.1
46	BU	27	VAL	2.1
45	DS	32	ALA	2.1
47	DF	170	ALA	2.1
1	AA	1226	C	2.1
43	DO	52	SER	2.1
47	DF	17	THR	2.1
28	BP	32	VAL	2.1
47	DF	147	ARG	2.1
8	AI	27	ILE	2.1
14	AQ	38	LYS	2.1
14	CQ	10	ARG	2.1
25	BC	80	LEU	2.1
42	DN	70	THR	2.1
47	BF	73	VAL	2.1
50	DT	50	LEU	2.1
29	DE	86	ALA	2.1
11	CL	81	ILE	2.1
44	DQ	64	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
38	DM	66	ARG	2.1
6	AG	42	VAL	2.1
7	AH	71	VAL	2.1
29	DE	189	THR	2.1
37	BL	125	LEU	2.1
39	BX	14	LEU	2.1
40	DH	142	VAL	2.1
25	BC	128	THR	2.1
41	BJ	140	LEU	2.1
23	DB	747	U	2.1
45	BS	75	PHE	2.1
25	BC	206	LYS	2.1
48	BG	43	LYS	2.1
25	BC	170	TYR	2.1
27	BK	32	TYR	2.1
44	DQ	63	ARG	2.1
39	DX	57	LEU	2.1
48	BG	89	VAL	2.1
48	BG	155	PRO	2.1
2	CC	141	MET	2.1
38	BM	113	ALA	2.1
44	BQ	102	LYS	2.1
24	DI	100	ILE	2.1
4	CE	72	ASN	2.1
27	DK	52	VAL	2.1
29	DE	3	LEU	2.1
41	BJ	25	LEU	2.1
51	BZ	6	GLN	2.1
3	CD	94	GLU	2.1
37	DL	114	GLY	2.1
38	BM	76	LYS	2.0
11	CL	16	ALA	2.0
29	DE	45	ALA	2.0
9	CJ	5	ARG	2.0
45	DS	3	THR	2.0
42	BN	79	LEU	2.0
35	BV	73	LYS	2.0
27	BK	38	ILE	2.0
38	DM	113	ALA	2.0
26	BD	12	THR	2.0
26	DD	83	ARG	2.0
4	CE	35	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
34	D3	56	LEU	2.0
7	AH	81	GLY	2.0
27	BK	96	GLY	2.0
7	AH	36	ALA	2.0
23	DB	690	G	2.0
7	CH	35	ILE	2.0
21	AN	55	SER	2.0
42	DN	29	VAL	2.0
47	DF	98	PHE	2.0
49	DR	95	ASP	2.0
3	AD	159	GLU	2.0
26	BD	147	GLY	2.0
7	CH	92	PRO	2.0
18	AB	99	MET	2.0
18	CB	22	TRP	2.0
50	DT	24	MET	2.0
35	DV	65	VAL	2.0
48	DG	112	VAL	2.0
47	DF	13	LYS	2.0
2	AC	205	GLU	2.0
2	CC	187	GLU	2.0
25	BC	120	ASP	2.0
15	AR	67	LEU	2.0
16	CS	12	LEU	2.0
25	BC	7	PRO	2.0
29	BE	156	ASN	2.0
25	DC	35	LYS	2.0
29	DE	64	GLY	2.0
35	BV	94	ALA	2.0
47	BF	44	ALA	2.0
52	BW	62	ALA	2.0
6	AG	22	LEU	2.0
9	CJ	65	TYR	2.0
24	DI	7	TYR	2.0
24	DI	81	LYS	2.0
25	DC	201	LEU	2.0
35	DV	61	LEU	2.0
26	BD	145	SER	2.0
45	DS	101	SER	2.0
23	BB	2310	C	2.0
49	DR	46	GLU	2.0
27	BK	22	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
46	BU	14	THR	2.0
47	DF	59	ILE	2.0
12	AM	80	MET	2.0
27	DK	62	VAL	2.0
28	BP	80	VAL	2.0
47	BF	37	MET	2.0
47	DF	167	ALA	2.0
18	AB	212	TYR	2.0
50	DT	84	TYR	2.0
43	DO	42	PRO	2.0
1	AA	1493	A	2.0
34	D3	7	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
53	MG	AA	1608	1/1	0.56	0.13	123,123,123,123	0
53	MG	BB	3042	1/1	0.60	0.23	92,92,92,92	0
53	MG	BB	3097	1/1	0.60	0.07	113,113,113,113	0
53	MG	AA	1622	1/1	0.70	0.34	130,130,130,130	0
53	MG	AA	1624	1/1	0.70	0.24	83,83,83,83	0
53	MG	DB	3052	1/1	0.77	0.29	105,105,105,105	0
53	MG	DB	3058	1/1	0.77	0.09	139,139,139,139	0
53	MG	AA	1635	1/1	0.77	0.15	80,80,80,80	0
53	MG	AA	1620	1/1	0.79	0.06	60,60,60,60	0
53	MG	DB	3094	1/1	0.79	0.14	100,100,100,100	0
53	MG	BB	3033	1/1	0.79	0.34	102,102,102,102	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	DB	3059	1/1	0.81	0.11	99,99,99,99	0
53	MG	CA	1646	1/1	0.82	0.11	139,139,139,139	0
53	MG	AA	1625	1/1	0.82	0.39	79,79,79,79	1
53	MG	AA	1623	1/1	0.83	0.37	33,33,33,33	1
53	MG	CA	1650	1/1	0.83	0.10	105,105,105,105	0
53	MG	CA	1612	1/1	0.84	0.28	93,93,93,93	0
53	MG	CA	1618	1/1	0.84	0.13	73,73,73,73	0
53	MG	BB	3049	1/1	0.84	0.19	67,67,67,67	0
53	MG	AA	1614	1/1	0.84	0.18	119,119,119,119	0
53	MG	CA	1651	1/1	0.85	0.08	101,101,101,101	0
53	MG	AA	1647	1/1	0.85	0.81	113,113,113,113	0
53	MG	CA	1640	1/1	0.85	0.11	62,62,62,62	0
53	MG	AA	1646	1/1	0.85	0.10	84,84,84,84	0
53	MG	CA	1644	1/1	0.86	0.14	57,57,57,57	0
53	MG	AA	1652	1/1	0.86	0.08	84,84,84,84	0
53	MG	DB	3029	1/1	0.86	0.70	87,87,87,87	0
53	MG	BB	3037	1/1	0.87	0.13	45,45,45,45	0
53	MG	CA	1613	1/1	0.88	0.49	126,126,126,126	0
53	MG	BB	3080	1/1	0.88	0.15	53,53,53,53	0
53	MG	BB	3093	1/1	0.88	0.11	108,108,108,108	0
53	MG	DB	3013	1/1	0.88	0.08	52,52,52,52	0
53	MG	BB	3035	1/1	0.88	0.07	13,13,13,13	0
53	MG	DB	3045	1/1	0.88	0.11	110,110,110,110	0
53	MG	CA	1648	1/1	0.88	0.07	104,104,104,104	0
53	MG	CA	1606	1/1	0.88	0.14	106,106,106,106	0
53	MG	AA	1626	1/1	0.88	0.20	36,36,36,36	1
53	MG	DB	3066	1/1	0.89	0.33	65,65,65,65	0
53	MG	AA	1655	1/1	0.89	0.14	88,88,88,88	0
53	MG	BB	3009	1/1	0.89	0.10	98,98,98,98	0
53	MG	AA	1606	1/1	0.89	0.07	82,82,82,82	0
53	MG	AA	1637	1/1	0.89	0.32	99,99,99,99	0
53	MG	AA	1617	1/1	0.89	0.09	112,112,112,112	0
53	MG	DB	3063	1/1	0.90	0.17	72,72,72,72	0
53	MG	BB	3010	1/1	0.90	0.16	53,53,53,53	0
53	MG	CA	1649	1/1	0.90	0.27	123,123,123,123	0
53	MG	BB	3068	1/1	0.90	0.19	13,13,13,13	0
53	MG	AA	1642	1/1	0.90	0.14	49,49,49,49	0
53	MG	AA	1649	1/1	0.90	0.04	89,89,89,89	0
53	MG	CA	1616	1/1	0.90	0.33	58,58,58,58	1
53	MG	DB	3032	1/1	0.90	0.16	73,73,73,73	0
53	MG	DB	3034	1/1	0.91	0.08	57,57,57,57	0
53	MG	BB	3051	1/1	0.91	0.12	107,107,107,107	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	CA	1629	1/1	0.91	0.17	67,67,67,67	0
53	MG	AA	1612	1/1	0.91	0.08	61,61,61,61	0
53	MG	AA	1627	1/1	0.91	0.07	63,63,63,63	0
53	MG	CA	1634	1/1	0.91	0.11	74,74,74,74	0
53	MG	AA	1657	1/1	0.91	0.34	91,91,91,91	0
53	MG	CA	1608	1/1	0.91	0.05	76,76,76,76	0
53	MG	BB	3046	1/1	0.91	0.10	89,89,89,89	0
53	MG	CA	1654	1/1	0.91	0.22	105,105,105,105	0
54	SCM	CA	1659	23/23	0.91	0.16	18,18,18,18	0
53	MG	AA	1656	1/1	0.92	0.14	87,87,87,87	0
53	MG	BB	3081	1/1	0.92	0.10	30,30,30,30	0
53	MG	AA	1605	1/1	0.92	0.10	50,50,50,50	0
53	MG	CA	1609	1/1	0.92	0.10	85,85,85,85	0
53	MG	DB	3035	1/1	0.92	0.22	79,79,79,79	0
53	MG	BB	3014	1/1	0.92	0.07	37,37,37,37	0
53	MG	BB	3001	1/1	0.92	0.19	5,5,5,5	0
53	MG	DN	201	1/1	0.92	0.51	145,145,145,145	0
53	MG	BB	3043	1/1	0.92	0.12	107,107,107,107	0
53	MG	CE	201	1/1	0.92	0.08	109,109,109,109	0
53	MG	CA	1628	1/1	0.92	0.11	52,52,52,52	0
53	MG	BB	3004	1/1	0.92	0.05	80,80,80,80	0
53	MG	DB	3104	1/1	0.92	0.12	29,29,29,29	0
53	MG	CA	1627	1/1	0.92	0.07	27,27,27,27	0
53	MG	DB	3050	1/1	0.92	0.11	80,80,80,80	0
53	MG	AA	1633	1/1	0.92	0.05	65,65,65,65	0
53	MG	AA	1650	1/1	0.92	0.08	94,94,94,94	0
55	ZN	D4	101	1/1	0.93	0.05	62,62,62,62	0
53	MG	BB	3108	1/1	0.93	0.06	10,10,10,10	0
53	MG	CA	1645	1/1	0.93	0.10	55,55,55,55	0
53	MG	DB	3064	1/1	0.93	0.39	49,49,49,49	0
53	MG	CA	1617	1/1	0.93	0.18	88,88,88,88	0
53	MG	CA	1632	1/1	0.93	0.24	76,76,76,76	0
53	MG	DB	3005	1/1	0.93	0.17	25,25,25,25	0
53	MG	AA	1621	1/1	0.94	0.17	36,36,36,36	0
53	MG	BB	3064	1/1	0.94	0.19	78,78,78,78	0
53	MG	DB	3028	1/1	0.94	0.18	70,70,70,70	0
53	MG	DB	3067	1/1	0.94	0.06	5,5,5,5	0
53	MG	DB	3022	1/1	0.94	0.06	11,11,11,11	0
53	MG	DB	3109	1/1	0.94	0.04	9,9,9,9	0
53	MG	AA	1645	1/1	0.94	0.20	70,70,70,70	0
53	MG	BB	3020	1/1	0.94	0.31	6,6,6,6	0
53	MG	CA	1625	1/1	0.94	0.05	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
55	ZN	B4	101	1/1	0.94	0.07	67,67,67,67	0
53	MG	BB	3071	1/1	0.94	0.18	68,68,68,68	0
53	MG	BB	3008	1/1	0.94	0.12	82,82,82,82	0
53	MG	CA	1637	1/1	0.94	0.17	98,98,98,98	0
53	MG	BB	3079	1/1	0.94	0.07	38,38,38,38	0
53	MG	AA	1628	1/1	0.94	0.18	70,70,70,70	0
53	MG	AA	1602	1/1	0.94	0.12	85,85,85,85	0
53	MG	DB	3091	1/1	0.94	0.11	90,90,90,90	0
53	MG	AA	1658	1/1	0.94	0.04	120,120,120,120	0
53	MG	BB	3005	1/1	0.94	0.11	5,5,5,5	0
53	MG	DB	3016	1/1	0.94	0.08	5,5,5,5	0
53	MG	DB	3071	1/1	0.94	0.17	57,57,57,57	0
53	MG	BB	3092	1/1	0.95	0.09	46,46,46,46	0
53	MG	CA	1615	1/1	0.95	0.06	13,13,13,13	0
53	MG	BB	3104	1/1	0.95	0.17	26,26,26,26	0
53	MG	DB	3096	1/1	0.95	0.11	26,26,26,26	0
53	MG	DB	3110	1/1	0.95	0.12	84,84,84,84	0
53	MG	BB	3003	1/1	0.95	0.06	53,53,53,53	0
53	MG	DB	3060	1/1	0.95	0.12	83,83,83,83	0
53	MG	AA	1618	1/1	0.95	0.08	38,38,38,38	0
53	MG	CA	1604	1/1	0.95	0.10	20,20,20,20	0
53	MG	BB	3015	1/1	0.95	0.10	25,25,25,25	0
53	MG	BB	3090	1/1	0.95	0.10	88,88,88,88	0
53	MG	BB	3099	1/1	0.95	0.06	41,41,41,41	0
53	MG	DB	3065	1/1	0.95	0.07	128,128,128,128	0
53	MG	BB	3025	1/1	0.95	0.15	30,30,30,30	0
53	MG	DB	3108	1/1	0.95	0.10	43,43,43,43	0
53	MG	BB	3078	1/1	0.95	0.19	27,27,27,27	0
53	MG	BB	3023	1/1	0.95	0.10	5,5,5,5	0
53	MG	DB	3003	1/1	0.95	0.05	30,30,30,30	0
53	MG	DB	3051	1/1	0.95	0.11	75,75,75,75	0
53	MG	DB	3018	1/1	0.95	0.09	7,7,7,7	0
53	MG	CA	1631	1/1	0.95	0.04	41,41,41,41	0
53	MG	AA	1654	1/1	0.95	0.07	67,67,67,67	0
53	MG	DB	3053	1/1	0.95	0.07	65,65,65,65	0
53	MG	CA	1605	1/1	0.95	0.05	12,12,12,12	0
53	MG	BB	3069	1/1	0.95	0.07	5,5,5,5	0
53	MG	BB	3017	1/1	0.96	0.13	50,50,50,50	0
53	MG	BB	3034	1/1	0.96	0.12	86,86,86,86	0
53	MG	DB	3089	1/1	0.96	0.13	79,79,79,79	0
53	MG	AA	1613	1/1	0.96	0.04	58,58,58,58	0
53	MG	BB	3054	1/1	0.96	0.13	49,49,49,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	BB	3006	1/1	0.96	0.15	61,61,61,61	0
53	MG	BB	3088	1/1	0.96	0.06	11,11,11,11	0
53	MG	DB	3055	1/1	0.96	0.16	44,44,44,44	0
53	MG	AA	1603	1/1	0.96	0.10	57,57,57,57	0
53	MG	BB	3070	1/1	0.96	0.05	74,74,74,74	0
53	MG	BB	3110	1/1	0.96	0.13	80,80,80,80	0
53	MG	AA	1632	1/1	0.96	0.21	96,96,96,96	0
53	MG	BB	3021	1/1	0.96	0.13	52,52,52,52	0
53	MG	BB	3032	1/1	0.96	0.07	22,22,22,22	0
53	MG	DB	3041	1/1	0.96	0.08	40,40,40,40	0
53	MG	DB	3027	1/1	0.96	0.07	13,13,13,13	0
53	MG	BB	3036	1/1	0.96	0.20	51,51,51,51	0
53	MG	AA	1615	1/1	0.96	0.35	96,96,96,96	0
53	MG	BB	3076	1/1	0.96	0.12	35,35,35,35	0
53	MG	BB	3038	1/1	0.96	0.08	92,92,92,92	0
53	MG	DB	3100	1/1	0.96	0.17	19,19,19,19	0
53	MG	CA	1623	1/1	0.96	0.10	11,11,11,11	0
53	MG	DB	3103	1/1	0.96	0.06	36,36,36,36	0
53	MG	BB	3024	1/1	0.96	0.09	55,55,55,55	0
53	MG	AA	1601	1/1	0.96	0.09	36,36,36,36	0
53	MG	CA	1658	1/1	0.96	0.08	37,37,37,37	0
53	MG	DB	3001	1/1	0.96	0.06	5,5,5,5	0
53	MG	BB	3052	1/1	0.96	0.10	71,71,71,71	0
53	MG	BB	3057	1/1	0.96	0.16	53,53,53,53	0
53	MG	CA	1641	1/1	0.96	0.10	42,42,42,42	0
53	MG	CA	1633	1/1	0.96	0.06	23,23,23,23	0
53	MG	AA	1619	1/1	0.97	0.04	111,111,111,111	0
53	MG	DB	3069	1/1	0.97	0.23	64,64,64,64	0
53	MG	BB	3067	1/1	0.97	0.17	63,63,63,63	0
53	MG	BB	3031	1/1	0.97	0.18	60,60,60,60	0
53	MG	BB	3065	1/1	0.97	0.06	32,32,32,32	0
53	MG	AA	1638	1/1	0.97	0.07	51,51,51,51	0
53	MG	DB	3030	1/1	0.97	0.19	47,47,47,47	0
53	MG	CA	1652	1/1	0.97	0.07	49,49,49,49	0
53	MG	DB	3057	1/1	0.97	0.05	48,48,48,48	0
53	MG	AA	1659	1/1	0.97	0.39	108,108,108,108	0
53	MG	CA	1653	1/1	0.97	0.07	48,48,48,48	0
53	MG	CA	1638	1/1	0.97	0.10	86,86,86,86	0
53	MG	DB	3078	1/1	0.97	0.04	25,25,25,25	0
53	MG	AA	1616	1/1	0.97	0.09	5,5,5,5	0
53	MG	BB	3102	1/1	0.97	0.12	43,43,43,43	0
53	MG	DB	3082	1/1	0.97	0.11	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	BB	3096	1/1	0.97	0.07	58,58,58,58	0
53	MG	BB	3084	1/1	0.97	0.06	32,32,32,32	0
53	MG	BB	3100	1/1	0.97	0.20	116,116,116,116	0
53	MG	DB	3010	1/1	0.97	0.14	5,5,5,5	0
53	MG	BB	3053	1/1	0.97	0.10	60,60,60,60	0
53	MG	DB	3031	1/1	0.97	0.09	46,46,46,46	0
53	MG	BB	3075	1/1	0.97	0.09	40,40,40,40	0
54	SCM	AA	1661	23/23	0.97	0.10	13,13,13,13	0
53	MG	BB	3082	1/1	0.97	0.15	18,18,18,18	0
53	MG	DB	3007	1/1	0.97	0.12	62,62,62,62	0
53	MG	DB	3054	1/1	0.97	0.15	21,21,21,21	0
53	MG	BB	3095	1/1	0.97	0.07	62,62,62,62	0
53	MG	CA	1607	1/1	0.97	0.06	5,5,5,5	0
53	MG	BB	3039	1/1	0.97	0.15	41,41,41,41	0
53	MG	DB	3079	1/1	0.97	0.06	7,7,7,7	0
53	MG	CA	1647	1/1	0.97	0.06	75,75,75,75	0
53	MG	AA	1631	1/1	0.97	0.12	5,5,5,5	0
53	MG	BB	3013	1/1	0.97	0.09	45,45,45,45	0
53	MG	AA	1629	1/1	0.97	0.06	12,12,12,12	0
53	MG	AA	1651	1/1	0.97	0.06	35,35,35,35	0
53	MG	DB	3106	1/1	0.97	0.05	61,61,61,61	0
53	MG	BB	3027	1/1	0.97	0.06	50,50,50,50	0
53	MG	AA	1660	1/1	0.97	0.03	56,56,56,56	0
53	MG	BB	3030	1/1	0.97	0.07	51,51,51,51	0
53	MG	DB	3046	1/1	0.98	0.04	23,23,23,23	0
53	MG	BB	3077	1/1	0.98	0.11	53,53,53,53	0
53	MG	DB	3061	1/1	0.98	0.04	66,66,66,66	0
53	MG	BB	3007	1/1	0.98	0.14	68,68,68,68	0
53	MG	BB	3062	1/1	0.98	0.08	7,7,7,7	0
53	MG	AA	1639	1/1	0.98	0.04	93,93,93,93	0
53	MG	DB	3004	1/1	0.98	0.08	8,8,8,8	0
53	MG	DB	3043	1/1	0.98	0.06	7,7,7,7	0
53	MG	AA	1636	1/1	0.98	0.04	64,64,64,64	0
53	MG	BB	3002	1/1	0.98	0.06	23,23,23,23	0
53	MG	DB	3017	1/1	0.98	0.23	5,5,5,5	0
53	MG	CA	1614	1/1	0.98	0.15	83,83,83,83	0
53	MG	CA	1635	1/1	0.98	0.06	30,30,30,30	0
53	MG	BB	3026	1/1	0.98	0.10	65,65,65,65	0
53	MG	DB	3047	1/1	0.98	0.15	43,43,43,43	0
53	MG	BB	3098	1/1	0.98	0.15	14,14,14,14	0
53	MG	DB	3048	1/1	0.98	0.11	38,38,38,38	0
53	MG	DB	3085	1/1	0.98	0.05	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	DB	3099	1/1	0.98	0.13	58,58,58,58	0
53	MG	BB	3089	1/1	0.98	0.12	72,72,72,72	0
53	MG	DB	3008	1/1	0.98	0.06	13,13,13,13	0
53	MG	BB	3087	1/1	0.98	0.25	80,80,80,80	0
53	MG	DB	3033	1/1	0.98	0.04	11,11,11,11	0
53	MG	BB	3056	1/1	0.98	0.03	5,5,5,5	0
53	MG	BB	3019	1/1	0.98	0.12	21,21,21,21	0
53	MG	DB	3092	1/1	0.98	0.15	10,10,10,10	0
53	MG	CA	1610	1/1	0.98	0.14	5,5,5,5	0
53	MG	DB	3107	1/1	0.98	0.04	5,5,5,5	0
53	MG	DB	3077	1/1	0.98	0.13	45,45,45,45	0
53	MG	CA	1619	1/1	0.98	0.09	36,36,36,36	0
53	MG	BB	3086	1/1	0.98	0.10	5,5,5,5	0
53	MG	BB	3044	1/1	0.98	0.09	45,45,45,45	0
53	MG	DB	3019	1/1	0.98	0.06	5,5,5,5	0
53	MG	AA	1609	1/1	0.98	0.06	40,40,40,40	0
53	MG	DB	3088	1/1	0.98	0.17	87,87,87,87	0
53	MG	BB	3103	1/1	0.98	0.08	11,11,11,11	0
53	MG	DB	3062	1/1	0.98	0.23	58,58,58,58	0
53	MG	BB	3061	1/1	0.98	0.08	24,24,24,24	0
53	MG	CA	1639	1/1	0.98	0.05	35,35,35,35	0
53	MG	DB	3074	1/1	0.98	0.07	26,26,26,26	0
53	MG	AA	1630	1/1	0.98	0.04	88,88,88,88	0
53	MG	BB	3074	1/1	0.98	0.06	10,10,10,10	0
53	MG	CA	1643	1/1	0.98	0.06	36,36,36,36	0
53	MG	DB	3090	1/1	0.98	0.09	14,14,14,14	0
53	MG	BB	3047	1/1	0.98	0.11	104,104,104,104	0
53	MG	DB	3036	1/1	0.98	0.04	87,87,87,87	0
53	MG	BB	3091	1/1	0.98	0.10	5,5,5,5	0
53	MG	DB	3087	1/1	0.98	0.08	5,5,5,5	0
53	MG	BB	3050	1/1	0.98	0.12	53,53,53,53	0
53	MG	BB	3048	1/1	0.98	0.04	12,12,12,12	0
53	MG	DB	3083	1/1	0.98	0.08	27,27,27,27	0
53	MG	BB	3018	1/1	0.98	0.12	39,39,39,39	0
53	MG	AA	1643	1/1	0.98	0.09	40,40,40,40	0
53	MG	DB	3056	1/1	0.98	0.06	5,5,5,5	0
53	MG	CA	1601	1/1	0.98	0.06	5,5,5,5	0
53	MG	BB	3040	1/1	0.98	0.08	12,12,12,12	0
53	MG	BB	3045	1/1	0.98	0.14	72,72,72,72	0
53	MG	DB	3070	1/1	0.98	0.06	33,33,33,33	0
53	MG	DB	3020	1/1	0.98	0.09	5,5,5,5	0
53	MG	DB	3095	1/1	0.98	0.15	92,92,92,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	DB	3006	1/1	0.98	0.11	10,10,10,10	0
53	MG	DB	3086	1/1	0.98	0.12	18,18,18,18	0
53	MG	DB	3072	1/1	0.98	0.07	20,20,20,20	0
53	MG	DB	3039	1/1	0.98	0.08	62,62,62,62	0
53	MG	CA	1626	1/1	0.98	0.08	8,8,8,8	0
53	MG	BB	3094	1/1	0.98	0.06	36,36,36,36	0
53	MG	AA	1640	1/1	0.98	0.15	46,46,46,46	0
53	MG	DB	3044	1/1	0.98	0.05	18,18,18,18	0
53	MG	BB	3011	1/1	0.98	0.15	5,5,5,5	0
53	MG	BB	3022	1/1	0.98	0.12	32,32,32,32	0
53	MG	BB	3109	1/1	0.98	0.10	11,11,11,11	0
53	MG	AA	1644	1/1	0.98	0.05	75,75,75,75	0
53	MG	DB	3009	1/1	0.98	0.07	9,9,9,9	0
53	MG	BB	3106	1/1	0.98	0.12	10,10,10,10	0
53	MG	DB	3002	1/1	0.98	0.05	11,11,11,11	0
53	MG	DB	3093	1/1	0.98	0.05	59,59,59,59	0
53	MG	DB	3098	1/1	0.98	0.10	5,5,5,5	0
53	MG	DB	3075	1/1	0.98	0.11	44,44,44,44	0
53	MG	DB	3024	1/1	0.98	0.05	40,40,40,40	0
53	MG	AA	1653	1/1	0.98	0.09	21,21,21,21	0
53	MG	CA	1621	1/1	0.98	0.16	80,80,80,80	0
53	MG	AA	1641	1/1	0.99	0.02	32,32,32,32	0
53	MG	DB	3073	1/1	0.99	0.07	50,50,50,50	0
53	MG	DB	3015	1/1	0.99	0.06	42,42,42,42	0
53	MG	CA	1602	1/1	0.99	0.12	5,5,5,5	0
53	MG	CA	1611	1/1	0.99	0.07	28,28,28,28	0
53	MG	BB	3072	1/1	0.99	0.10	35,35,35,35	0
53	MG	CA	1657	1/1	0.99	0.05	73,73,73,73	0
53	MG	DB	3076	1/1	0.99	0.10	33,33,33,33	0
53	MG	BB	3083	1/1	0.99	0.14	30,30,30,30	0
53	MG	DB	3026	1/1	0.99	0.23	45,45,45,45	0
53	MG	BB	3012	1/1	0.99	0.06	67,67,67,67	0
53	MG	DB	3023	1/1	0.99	0.04	69,69,69,69	0
53	MG	AA	1604	1/1	0.99	0.09	38,38,38,38	0
53	MG	BB	3059	1/1	0.99	0.13	10,10,10,10	0
53	MG	DB	3021	1/1	0.99	0.04	16,16,16,16	0
53	MG	BB	3073	1/1	0.99	0.10	44,44,44,44	0
53	MG	DB	3102	1/1	0.99	0.23	28,28,28,28	0
53	MG	CA	1630	1/1	0.99	0.10	37,37,37,37	0
53	MG	CA	1624	1/1	0.99	0.07	38,38,38,38	0
53	MG	DB	3097	1/1	0.99	0.17	41,41,41,41	0
53	MG	DB	3080	1/1	0.99	0.10	62,62,62,62	0

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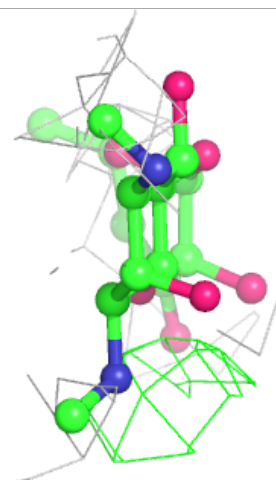
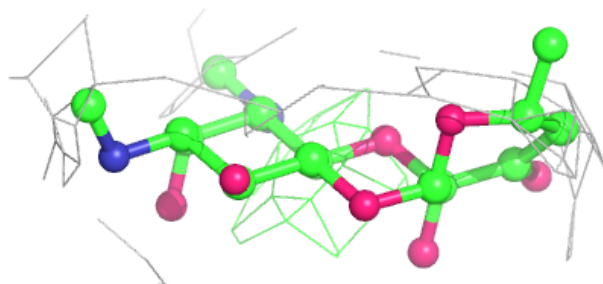
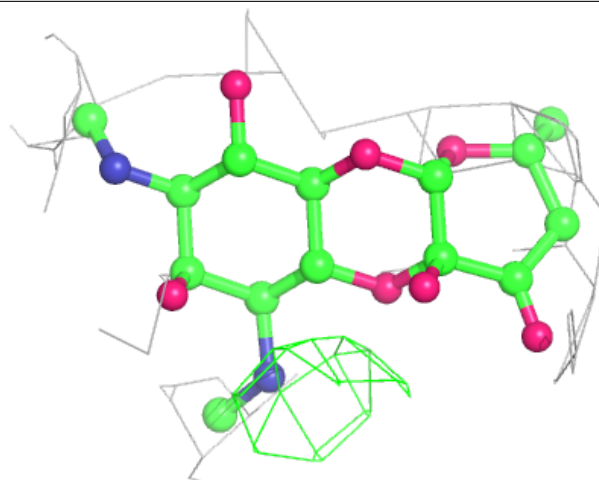
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
53	MG	BB	3063	1/1	0.99	0.05	33,33,33,33	0
53	MG	BB	3041	1/1	0.99	0.09	5,5,5,5	0
53	MG	DB	3068	1/1	0.99	0.15	5,5,5,5	0
53	MG	DB	3037	1/1	0.99	0.16	8,8,8,8	0
53	MG	BB	3105	1/1	0.99	0.18	5,5,5,5	0
53	MG	BB	3058	1/1	0.99	0.06	15,15,15,15	0
53	MG	DB	3040	1/1	0.99	0.05	5,5,5,5	0
53	MG	DB	3012	1/1	0.99	0.12	9,9,9,9	0
53	MG	CA	1655	1/1	0.99	0.03	32,32,32,32	0
53	MG	CA	1656	1/1	0.99	0.14	38,38,38,38	0
53	MG	DB	3081	1/1	0.99	0.06	41,41,41,41	0
53	MG	CA	1620	1/1	0.99	0.10	72,72,72,72	0
53	MG	BB	3107	1/1	0.99	0.21	46,46,46,46	0
53	MG	AA	1607	1/1	0.99	0.09	57,57,57,57	0
53	MG	BB	3085	1/1	0.99	0.07	34,34,34,34	0
53	MG	AA	1648	1/1	0.99	0.06	6,6,6,6	0
53	MG	DB	3042	1/1	0.99	0.04	37,37,37,37	0
53	MG	BB	3028	1/1	0.99	0.24	9,9,9,9	0
53	MG	BB	3029	1/1	0.99	0.05	5,5,5,5	0
53	MG	DB	3038	1/1	0.99	0.07	5,5,5,5	0
53	MG	BB	3016	1/1	0.99	0.04	55,55,55,55	0
53	MG	AA	1634	1/1	0.99	0.03	72,72,72,72	0
53	MG	BB	3066	1/1	0.99	0.04	5,5,5,5	0
53	MG	CA	1603	1/1	0.99	0.03	56,56,56,56	0
53	MG	BB	3060	1/1	0.99	0.17	30,30,30,30	0
53	MG	DB	3025	1/1	0.99	0.07	44,44,44,44	0
53	MG	AA	1611	1/1	0.99	0.05	43,43,43,43	0
53	MG	BB	3055	1/1	0.99	0.12	6,6,6,6	0
53	MG	DB	3105	1/1	0.99	0.04	40,40,40,40	0
53	MG	CA	1642	1/1	0.99	0.07	79,79,79,79	0
53	MG	AA	1610	1/1	0.99	0.06	78,78,78,78	0
53	MG	CA	1622	1/1	0.99	0.17	5,5,5,5	0
53	MG	DB	3084	1/1	0.99	0.09	29,29,29,29	0
53	MG	DB	3011	1/1	0.99	0.10	5,5,5,5	0
53	MG	DB	3101	1/1	0.99	0.10	26,26,26,26	0
53	MG	DB	3014	1/1	1.00	0.05	21,21,21,21	0
53	MG	BB	3101	1/1	1.00	0.05	19,19,19,19	0
53	MG	CA	1636	1/1	1.00	0.04	5,5,5,5	0
53	MG	DB	3049	1/1	1.00	0.14	42,42,42,42	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

orientation to approximate a three-dimensional view.

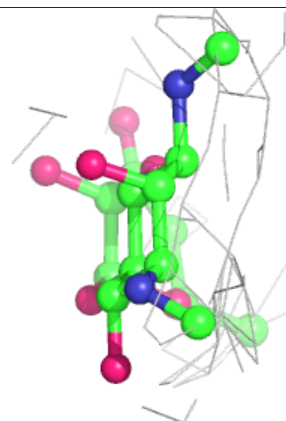
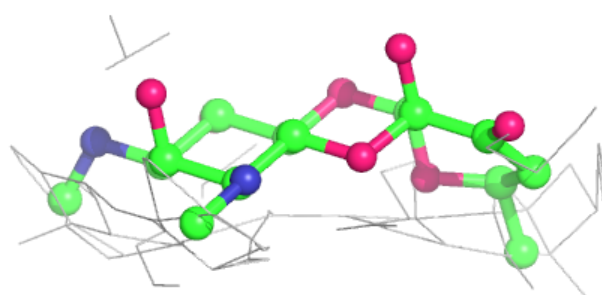
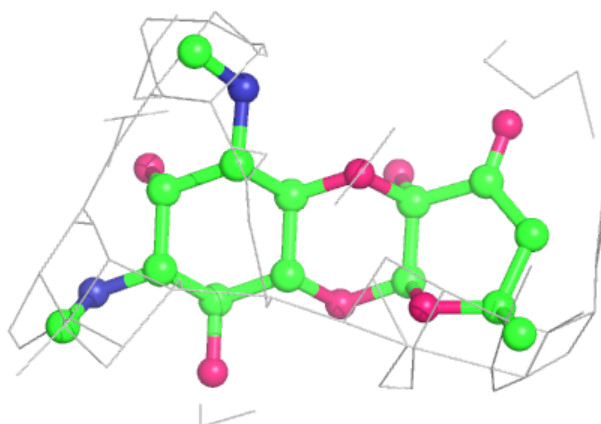
Electron density around SCM CA 1659:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SCM AA 1661:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.