



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 08:59 am GMT

PDB ID : 4V4H
Title : Crystal structure of the bacterial ribosome from Escherichia coli in complex with the antibiotic kasugamyin at 3.5Å resolution.
Authors : Schuwirth, B.S.; Vila-Sanjurjo, A.; Cate, J.H.D.
Deposited on : 2006-08-04
Resolution : 3.46 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28972

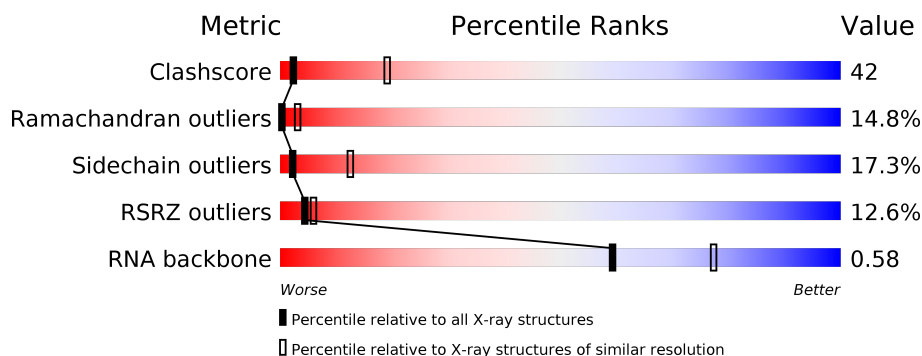
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.46 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))
Clashscore	112137	1040 (3.52-3.40)
Ramachandran outliers	110173	1009 (3.52-3.40)
Sidechain outliers	110143	1010 (3.52-3.40)
RSRZ outliers	101464	1017 (3.54-3.38)
RNA backbone	2435	1020 (4.02-2.86)

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. 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>23%</div> <div>63%</div> <div>13%</div> <div>.</div> </div>
1	CA	1542	<div> <div>22%</div> <div>65%</div> <div>11%</div> <div>.</div> </div>
2	AC	233	<div> <div>11%</div> <div>25%</div> <div>52%</div> <div>11%</div> <div>.</div> <div>12%</div> </div>
2	CC	233	<div> <div>5%</div> <div>30%</div> <div>45%</div> <div>12%</div> <div>.</div> <div>12%</div> </div>
3	AD	206	<div> <div>22%</div> <div>31%</div> <div>51%</div> <div>16%</div> <div>.</div> </div>
3	CD	206	<div> <div>27%</div> <div>56%</div> <div>16%</div> </div>

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Mol	Chain	Length	Quality of chain
4	AE	167	
4	CE	167	
5	AF	135	
5	CF	135	
6	AG	179	
6	CG	179	
7	AH	130	
7	CH	130	
8	AI	130	
8	CI	130	
9	AJ	103	
9	CJ	103	
10	AK	129	
10	CK	129	
11	AL	124	
11	CL	124	
12	AM	118	
12	CM	118	
13	AN	101	
13	CN	101	
14	AO	89	
14	CO	89	
15	AP	82	
15	CP	82	
16	AQ	84	

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Mol	Chain	Length	Quality of chain
16	CQ	84	
17	AR	75	
17	CR	75	
18	AS	92	
18	CS	92	
19	AT	87	
19	CT	87	
20	AB	241	
20	CB	241	
21	AU	71	
21	CU	71	
22	BA	120	
22	DA	120	
23	BB	2904	
23	DB	2904	
24	BV	94	
24	DV	94	
25	BC	273	
25	DC	273	
26	BD	209	
26	DD	209	
27	BE	201	
27	DE	201	
28	BF	179	
28	DF	179	

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Mol	Chain	Length	Quality of chain
29	BG	177	
29	DG	177	
30	BH	149	
30	DH	149	
31	BJ	142	
31	DJ	142	
32	BK	123	
32	DK	123	
33	BL	144	
33	DL	144	
34	BM	136	
34	DM	136	
35	BN	127	
35	DN	127	
36	BO	117	
36	DO	117	
37	BP	115	
37	DP	115	
38	BQ	118	
38	DQ	118	
39	BR	103	
39	DR	103	
40	BS	110	
40	DS	110	
41	BT	100	

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Mol	Chain	Length	Quality of chain
41	DT	100	
42	BU	104	
42	DU	104	
43	BW	85	
43	DW	85	
44	BX	63	
44	DX	63	
45	BY	59	
45	DY	59	
46	BZ	70	
46	DZ	70	
47	B0	57	
47	D0	57	
48	B1	55	
48	D1	55	
49	B2	46	
49	D2	46	
50	B3	65	
50	D3	65	
51	B4	38	
51	D4	38	
52	BI	142	
52	DI	142	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
53	KSG	AA	1601	-	-	-	X
53	KSG	CA	1601	-	-	-	X
54	MG	BB	3086	-	-	-	X
54	MG	BB	3100	-	-	-	X
54	MG	DB	3082	-	-	-	X

2 Entry composition

There are 55 unique types of molecules in this entry. The entry contains 284160 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	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 S14.

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

- Molecule 14 is a protein called 30S RIBOSOMAL PROTEIN S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AO	88	Total	C	N	O	S	0	0	0
			716	440	146	129	1			
14	CO	88	Total	C	N	O	S	0	0	0
			716	440	146	129	1			

- Molecule 15 is a protein called 30S RIBOSOMAL PROTEIN S16.

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

- Molecule 16 is a protein called 30S RIBOSOMAL PROTEIN S17.

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

- Molecule 17 is a protein called 30S RIBOSOMAL PROTEIN S18.

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

- Molecule 18 is a protein called 30S RIBOSOMAL PROTEIN S19.

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

- Molecule 19 is a protein called 30S RIBOSOMAL PROTEIN S20.

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

- Molecule 20 is a protein called 30S RIBOSOMAL PROTEIN S2.

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

- Molecule 21 is a protein called 30S RIBOSOMAL PROTEIN S21.

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

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

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			

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

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			

- Molecule 24 is a protein called 50S RIBOSOMAL PROTEIN L25.

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

- Molecule 25 is a protein called 50S RIBOSOMAL PROTEIN L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	BC	267	Total	C	N	O	S	0	0	0
			2053	1271	416	359	7			
25	DC	267	Total	C	N	O	S	0	0	0
			2053	1271	416	359	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			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
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 L4.

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

- Molecule 28 is a protein called 50S RIBOSOMAL PROTEIN L5.

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

- Molecule 29 is a protein called 50S RIBOSOMAL PROTEIN L6.

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

- Molecule 30 is a protein called 50S RIBOSOMAL PROTEIN L9.

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

- Molecule 31 is a protein called 50S RIBOSOMAL PROTEIN L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	BJ	140	Total	C	N	O	S	0	0	0
			1112	704	210	194	4			
31	DJ	140	Total	C	N	O	S	0	0	0
			1112	704	210	194	4			

- Molecule 32 is a protein called 50S RIBOSOMAL PROTEIN L14.

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

- Molecule 33 is a protein called 50S RIBOSOMAL PROTEIN L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	BL	144	Total	C	N	O	S	0	0	0
			1053	654	207	190	2			
33	DL	144	Total	C	N	O	S	0	0	0
			1053	654	207	190	2			

- Molecule 34 is a protein called 50S RIBOSOMAL PROTEIN L16.

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

- Molecule 35 is a protein called 50S RIBOSOMAL PROTEIN L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BN	127	Total	C	N	O	S	0	0	0
			1008	621	204	178	5			
35	DN	127	Total	C	N	O	S	0	0	0
			1008	621	204	178	5			

- Molecule 36 is a protein called 50S RIBOSOMAL PROTEIN L18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	BO	117	Total	C	N	O	S	0	0	0
			900	557	179	163	1			
36	DO	117	Total	C	N	O	S	0	0	0
			900	557	179	163	1			

- Molecule 37 is a protein called 50S RIBOSOMAL PROTEIN L19.

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

- Molecule 38 is a protein called 50S RIBOSOMAL PROTEIN L20.

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

- Molecule 39 is a protein called 50S RIBOSOMAL PROTEIN L21.

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

- Molecule 40 is a protein called 50S RIBOSOMAL PROTEIN L22.

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

- Molecule 41 is a protein called 50S RIBOSOMAL PROTEIN L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BT	99	Total	C	N	O	S	0	0	0
			777	491	145	139	2			
41	DT	99	Total	C	N	O	S	0	0	0
			777	491	145	139	2			

- Molecule 42 is a protein called 50S RIBOSOMAL PROTEIN L24.

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

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

- Molecule 43 is a protein called 50S RIBOSOMAL PROTEIN L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	BW	84	Total	C	N	O	S			
			634	391	129	113	1	0	0	0
43	DW	84	Total	C	N	O	S			
			634	391	129	113	1	0	0	0

- Molecule 44 is a protein called 50S RIBOSOMAL PROTEIN L29.

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

- Molecule 45 is a protein called 50S RIBOSOMAL PROTEIN L30.

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

- Molecule 46 is a protein called 50S RIBOSOMAL PROTEIN L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	BZ	70	Total	C	N	O	S			
			549	339	104	100	6	0	0	0
46	DZ	70	Total	C	N	O	S			
			549	339	104	100	6	0	0	0

- Molecule 47 is a protein called 50S RIBOSOMAL PROTEIN L32.

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

- Molecule 48 is a protein called 50S RIBOSOMAL PROTEIN L33.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
48	B1	54	Total	C	N	O	0	0	0
			441	284	81	76			
48	D1	54	Total	C	N	O	0	0	0
			441	284	81	76			

- Molecule 49 is a protein called 50S RIBOSOMAL PROTEIN L34.

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

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

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

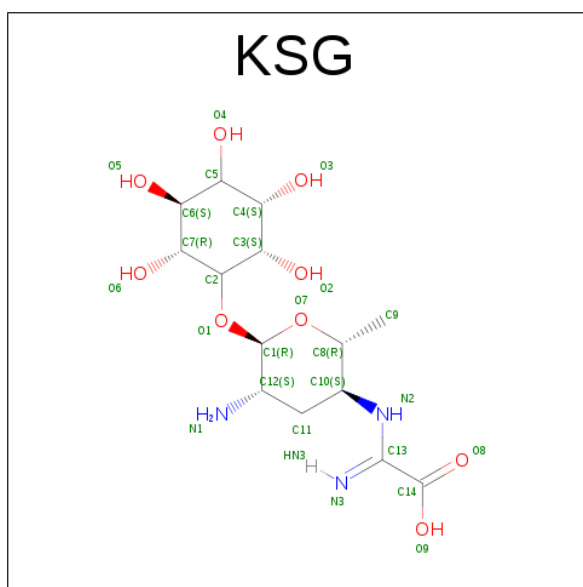
- Molecule 51 is a protein called 50S RIBOSOMAL PROTEIN L36.

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

- Molecule 52 is a protein called 50S RIBOSOMAL PROTEIN L11.

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

- Molecule 53 is (1S,2R,3S,4R,5S,6S)-2,3,4,5,6-PENTAHYDROXYCYCLOHEXYL 2-AMINO-4-{[CARBOXY(IMINO)METHYL]AMINO}-2,3,4,6-TETRADEOXY-ALPHA-D-ARABINO-HEXOPYRANOSIDE (three-letter code: KSG) (formula: C₁₄H₂₅N₃O₉).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
53	AA	1	Total	C	N	O	0	0
			26	14	3	9		
53	CA	1	Total	C	N	O	0	0
			26	14	3	9		

- Molecule 54 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	BB	110	Total	Mg	0	0
			110	110		
54	DE	1	Total	Mg	0	0
			1	1		
54	CA	62	Total	Mg	0	0
			62	62		
54	AA	60	Total	Mg	0	0
			60	60		
54	DN	1	Total	Mg	0	0
			1	1		
54	DB	109	Total	Mg	0	0
			109	109		

- Molecule 55 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
55	AA	289	Total	O	0	0
			289	289		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
55	AE	3	Total O 3 3	0	0
55	AK	2	Total O 2 2	0	0
55	AN	3	Total O 3 3	0	0
55	AP	2	Total O 2 2	0	0
55	AT	1	Total O 1 1	0	0
55	BB	497	Total O 497 497	0	0
55	BC	1	Total O 1 1	0	0
55	BE	5	Total O 5 5	0	0
55	BH	1	Total O 1 1	0	0
55	BL	2	Total O 2 2	0	0
55	BN	1	Total O 1 1	0	0
55	CA	293	Total O 293 293	0	0
55	CE	3	Total O 3 3	0	0
55	CK	1	Total O 1 1	0	0
55	CL	4	Total O 4 4	0	0
55	CN	3	Total O 3 3	0	0
55	CP	1	Total O 1 1	0	0
55	CT	3	Total O 3 3	0	0
55	DB	501	Total O 501 501	0	0
55	DC	1	Total O 1 1	0	0
55	DD	1	Total O 1 1	0	0

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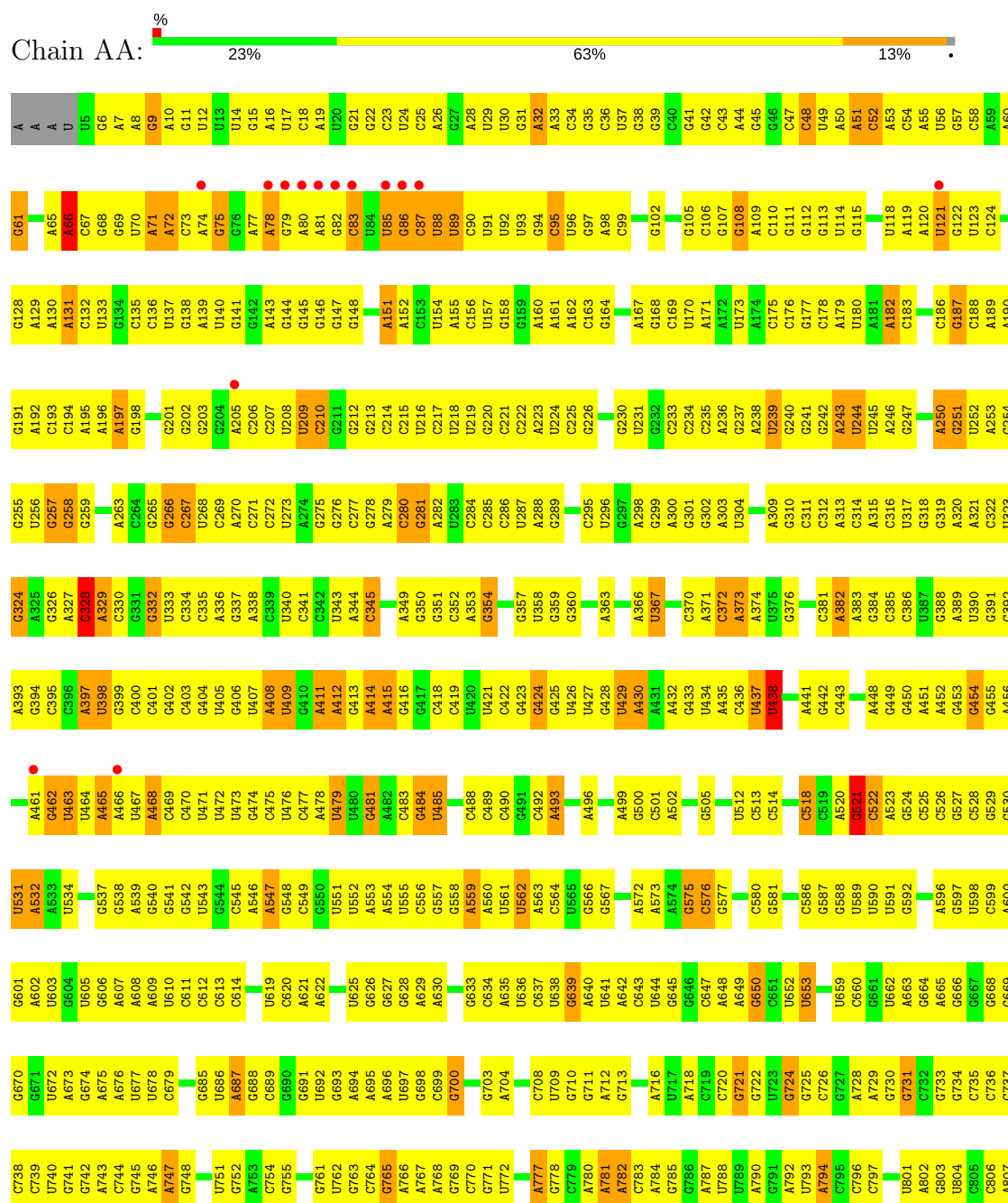
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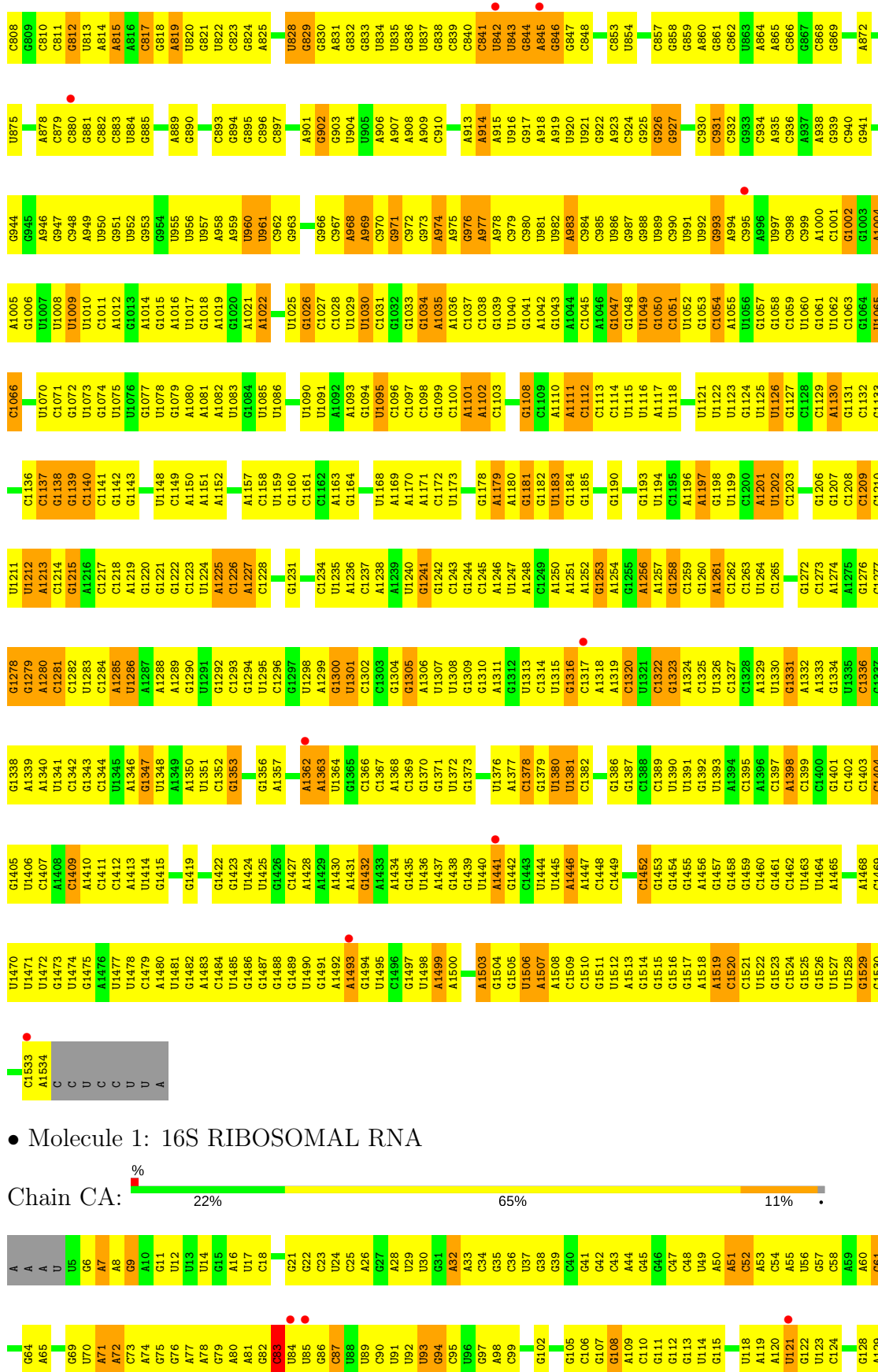
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
55	DE	3	Total 3	O 3	0	0
55	DL	1	Total 1	O 1	0	0
55	DN	2	Total 2	O 2	0	0
55	DT	1	Total 1	O 1	0	0
55	D2	2	Total 2	O 2	0	0

3 Residue-property plots

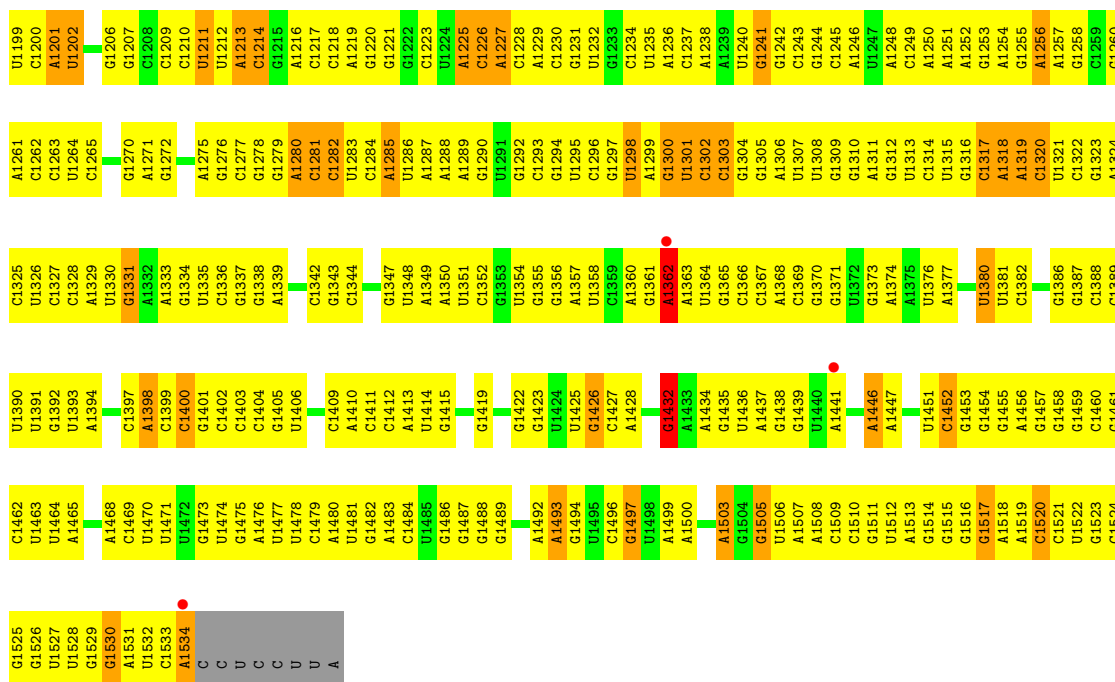
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

• Molecule 1: 16S RIBOSOMAL RNA

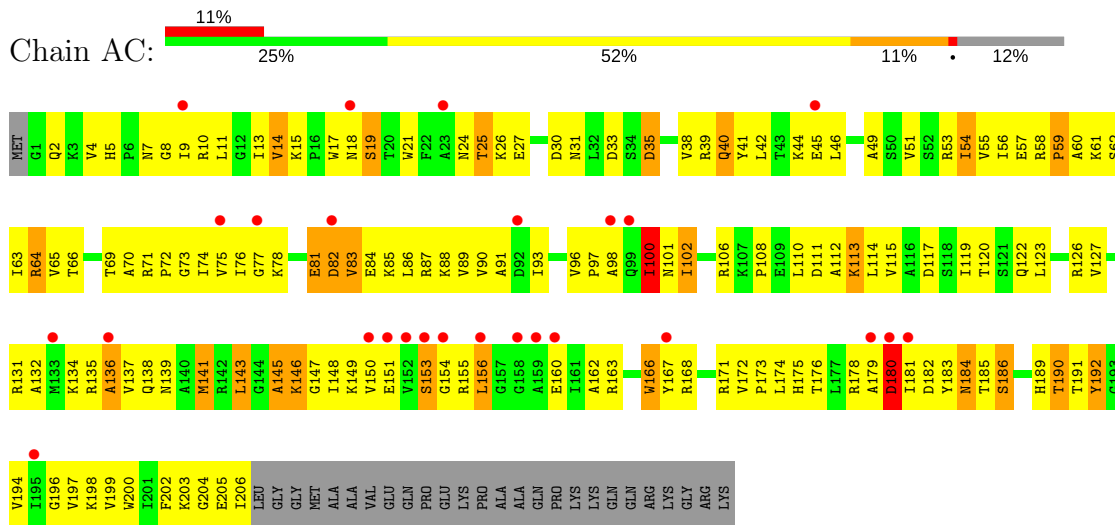




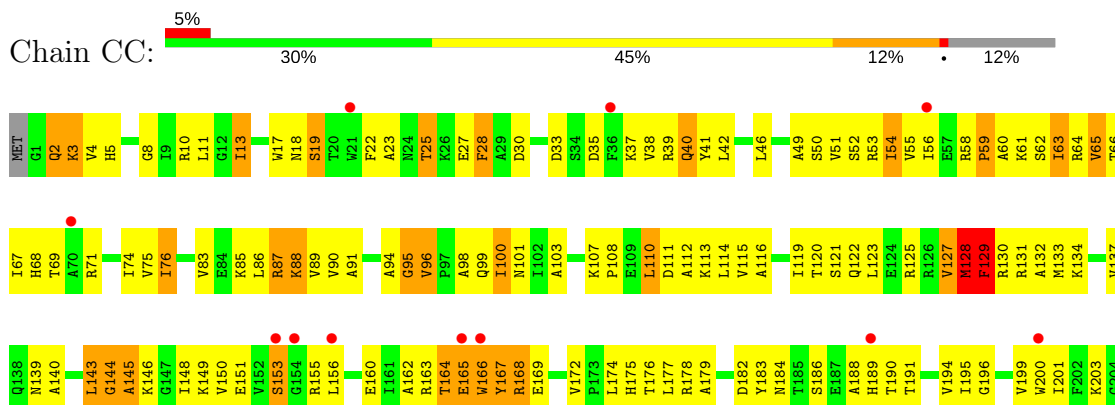
U1135	A1067	G1006	G945	U875	C808	C736	G669	U605	G537	U463	A397	A327	G259	C194	A130
C1136	G1068	U1007	A946	A878	G809	C737	U672	G606	G538	U464	U398	C328	A262	A196	A131
C1137	U1008	U1008	G947	C879	C810	C738	U673	A607	G539	A465	G399	A329	A263	A197	C132
G1138	C1070	U1009	C948	C880	C811	C739	A673	A608	A539	A466	C400	A330	C264	G198	U133
C1139	C1071	U1010	A949	C881	G812	U740	G674	A609	G540	U467	C401	G331	G265	G134	C135
C1140	G1072	G1011	U950	C882	U813	G741	A675	U610	G541	A468	G402	G332	G266	G202	G138
C1141	U1073	A1012	G951	C883	A814	G742	A676	C611	G542	C469	C403	U333	C267	G203	A139
G1142	G1074	G1013	U952	C884	A815	A743	A677	C612	U543	C470	G404	U334	C268	G204	U140
G1143	U1075	G1015	G953	C885	A816	A744	U678	C613	G544	U471	U405	C335	U268	A205	C206
A1145	U1076	G1016	G954	C886	C817	G745	C679	C614	A546	U472	G406	A336	C269	C270	G141
A1146	G1077	A1017	U955	C887	G818	A746	U684	G617	A547	C473	U407	A338	C271	C207	G142
C1147	U1078	U1018	U956	C888	A819	A747	G685	G618	A548	C474	A408	C339	C272	U208	A143
C1148	G1079	G1019	U957	C889	U820	G748	G686	U619	C549	U476	G410	U340	U273	U209	G144
C1149	A1080	G1020	A953	C890	G821	A749	U687	C620	G550	C477	A411	C341	A274	C210	G145
A1150	A1082	A1021	U959	C891	U822	C750	A687	U621	G551	A478	A412	C342	G275	G211	G147
A1151	U1083	C823	U960	C892	C824	U751	G688	A621	U552	U479	G413	U343	G276	G212	G148
A1152	G1084	A1022	U961	C893	G825	G752	G689	A622	U553	U480	A414	A344	C277	G213	G149
G1153	U1085	U1023	C962	C894	A826	A753	G690	U625	A554	A481	A415	C345	G278	C214	A149
G1154	U1086	G1024	G963	C895	A827	C754	G691	U626	U555	A482	G416	C346	C279	C215	U150
A1155	U1025	U1025	A964	A900	U828	G755	U692	G627	U556	C483	C417	C347	C280	U216	A151
A1156	G1026	G1026	U965	A901	G829	G756	U693	G628	C556	C484	C418	A349	C281	U217	A152
G1157	C1027	C1027	G966	G902	G830	G761	A694	U629	G557	U485	C419	G350	A282	U218	C152
C1158	C1028	G1033	C967	G903	A831	U762	A695	A630	G558	U486	C420	G351	U283	U219	U154
U1159	U1029	U1030	A968	U904	G832	G763	A696	A630	U559	C488	U421	C352	C284	G220	A155
U1160	U1030	C1031	A969	U905	G833	C764	G697	A630	U561	C489	C422	A353	C285	C221	C156
C1161	C1032	C1032	C972	A906	U834	G765	G698	G633	U562	C490	G423	G354	C286	C222	U157
C1162	G1033	G1033	G973	A907	U835	A766	C699	C634	A563	C491	G424	G357	U287	A223	G158
A1163	U1034	G1034	A974	A908	G836	A767	G700	C635	U564	C492	A425	U358	C288	U224	C159
C1164	G1035	A1035	A975	C910	U837	A768	G701	U644	C576	C493	U426	G359	G289	C225	A160
U1165	U1036	U1036	G976	U911	G838	C770	A704	U638	G566	C496	U427	C360	C295	G226	A161
A1166	C1037	C1037	A977	C912	C840	G771	G705	G639	G567	U496	G428	C361	C296	G230	A162
A1167	U1038	C1038	A978	A913	C841	U772	A706	A640	U572	A499	U429	C362	G297	U231	C163
U1168	C1039	C979	C980	A914	U842	U773	G707	U641	A573	C500	A430	U367	G298	G232	G164
A1169	G1040	G1040	C981	A915	U843	A777	C708	A642	A574	C501	A431	C370	G299	C233	A167
A1170	U1041	G1041	U981	U916	G844	C778	G709	C643	A575	A502	G433	A371	A300	C234	G168
A1171	G106	A1042	U982	G917	A845	C779	G710	U644	C576	U505	U434	C372	G301	C235	C169
C1172	C1107	G1043	A983	A918	G846	A780	G711	G645	G577	C505	A435	A373	G302	A236	U170
U1173	U1044	A1044	C984	A919	G847	A781	A712	G646	C577	U512	U437	A374	A303	G237	U173
G1174	C1045	C1045	C985	U920	C848	A782	G713	C647	C580	C513	U439	U375	A304	A238	A174
G1175	U1046	A1046	U986	U921	C853	C783	A716	A648	C581	C514	U440	G376	A309	G240	C175
A1176	G1047	G1047	G987	G922	U854	A784	A649	A650	C586	C518	A441	C381	A310	G241	C176
G1177	U1048	G1048	G988	A923	U855	G785	G719	C651	G587	C520	G442	C382	G312	G242	G178
G1178	U1049	U1049	U989	C924	U856	G786	C720	U652	G588	C521	C443	A383	A313	A243	A179
A1179	C1119	G1050	C990	G925	C857	U788	G721	U653	U589	G522	G448	C384	G314	U244	U180
A1180	C1120	C1051	U991	G926	G858	U789	G722	G656	U590	A523	A449	C385	A315	A246	A181
G1184	U1122	G1053	G992	G927	G859	A792	U723	G657	U591	G524	G450	U387	C316	G247	A182
G1185	C1054	C1054	G993	G928	A860	U793	G724	U657	G592	C525	G451	G388	U317	A250	C183
G1186	A1055	A1055	A994	G929	G861	U794	G725	G658	U596	C526	A452	G389	G319	G251	C186
G1187	U1125	U1056	C995	C930	C862	A794	G726	U659	U597	C527	G453	U390	G320	U252	G187
A1188	U1126	U1057	A996	C931	U863	C797	G727	U660	G598	C528	G454	A321	A321	A253	C188
G1189	G1127	U1060	C998	C932	A864	C797	G728	C661	U599	C529	G455	C392	G322	G254	A189
U1190	C1128	C999	C999	G933	A865	U801	A729	U662	U598	C530	G456	A393	U323	G255	A190
A1191	A1000	G1061	A1000	C934	C866	U802	G730	A663	G599	G530	A457	C394	U324	U256	G191
C1192	A1130	U1062	C1001	A935	G867	A803	G731	A664	G601	U531	A458	G395	A325	G257	A192
G1193	G1131	C1063	G1002	G839	G868	C803	C732	A665	G602	A532	A461	C396	G326	G258	C193
A1196	C1132	U1064	A1003	C940	G869	U804	C733	G666	U603	U534	G462	C397			
A1197	G1133	U1065	A1004	C941	A872	C806	G734	G667	G604						
G1198	G1134	C1066	A1005			A807	C735	G668							



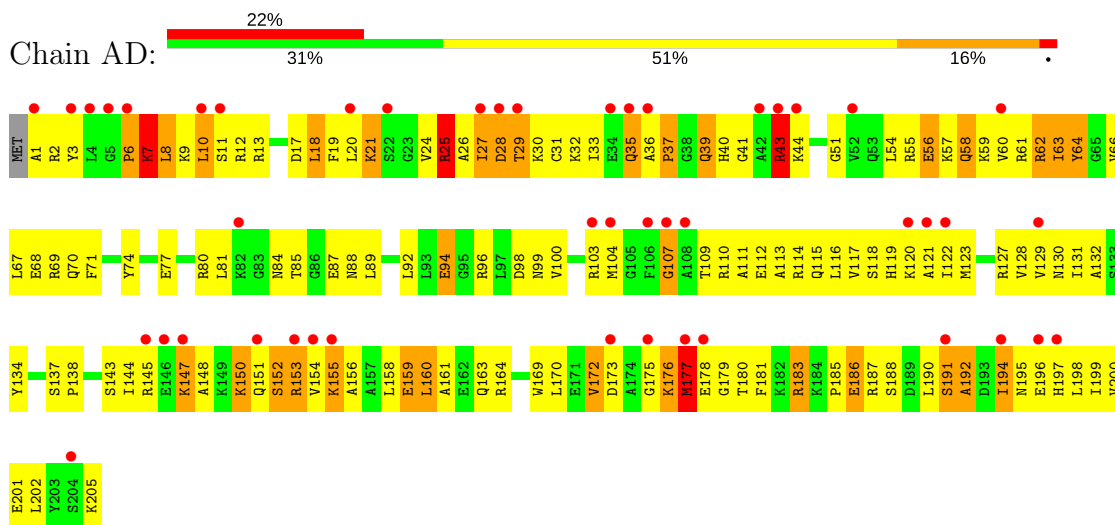
- Molecule 2: 30S RIBOSOMAL PROTEIN S3



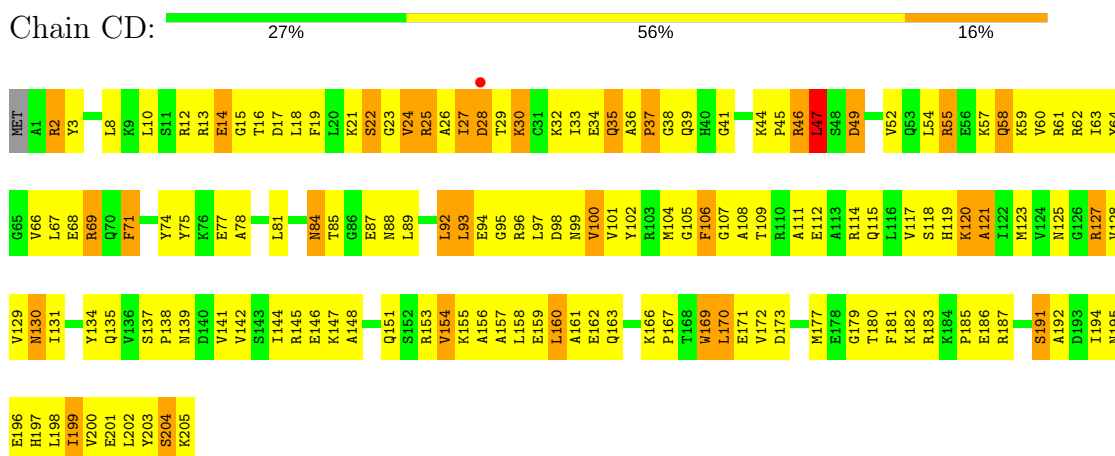
- Molecule 2: 30S RIBOSOMAL PROTEIN S3



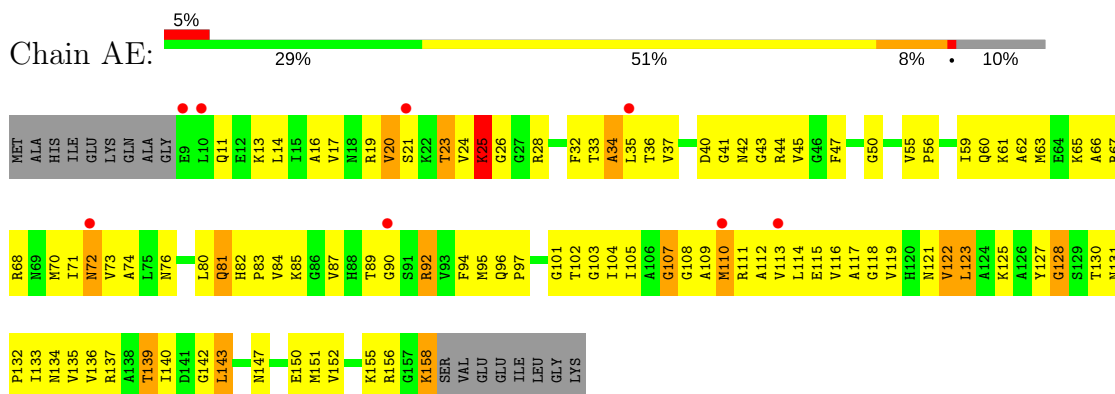
- Molecule 3: 30S RIBOSOMAL PROTEIN S4



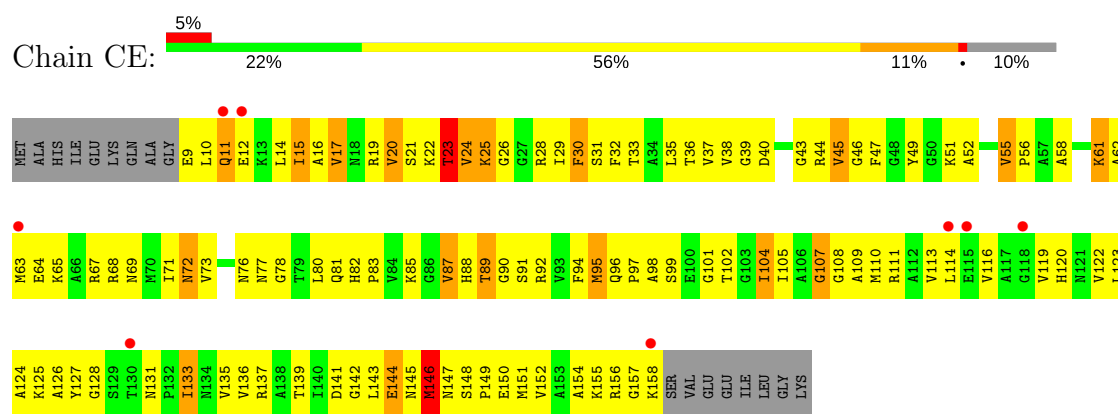
- Molecule 3: 30S RIBOSOMAL PROTEIN S4



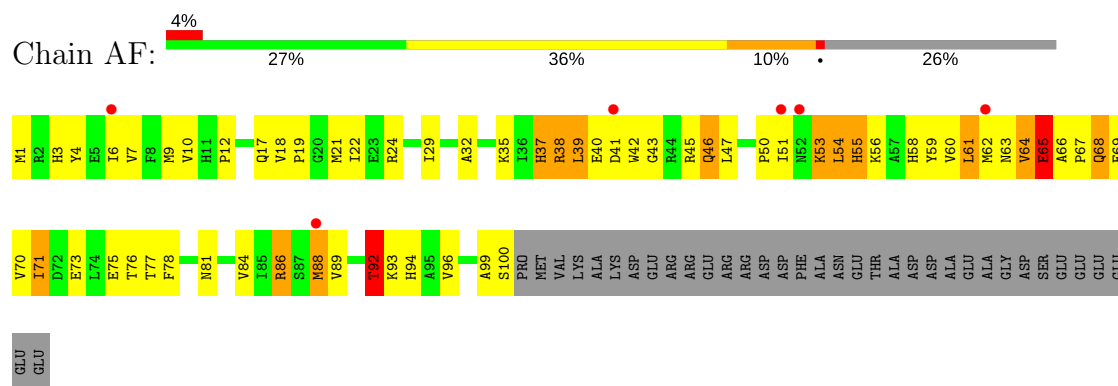
- Molecule 4: 30S RIBOSOMAL PROTEIN S5



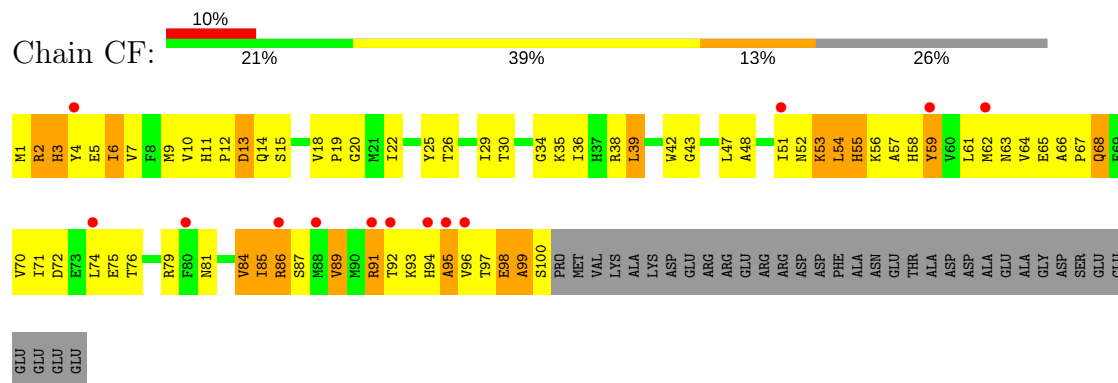
- Molecule 4: 30S RIBOSOMAL PROTEIN S5



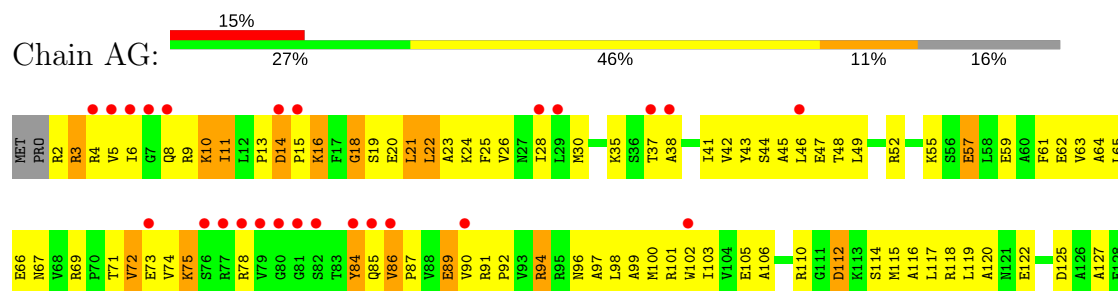
• Molecule 5: 30S RIBOSOMAL PROTEIN S6

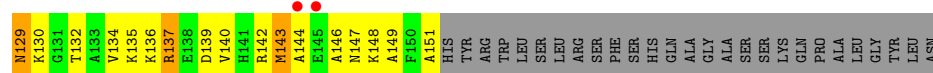


• Molecule 5: 30S RIBOSOMAL PROTEIN S6

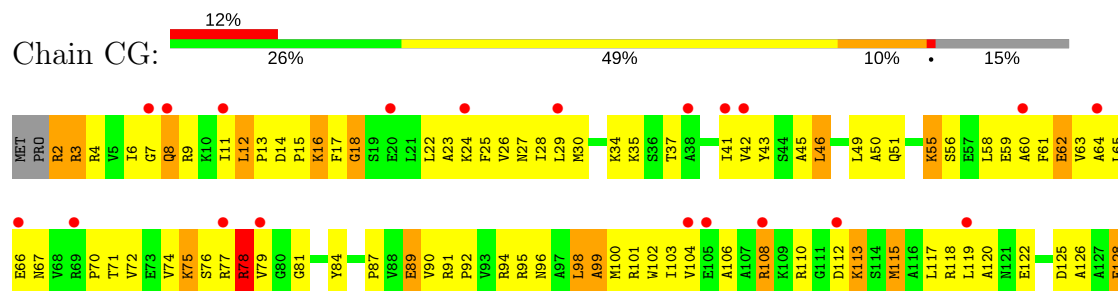


• Molecule 6: 30S RIBOSOMAL PROTEIN S7

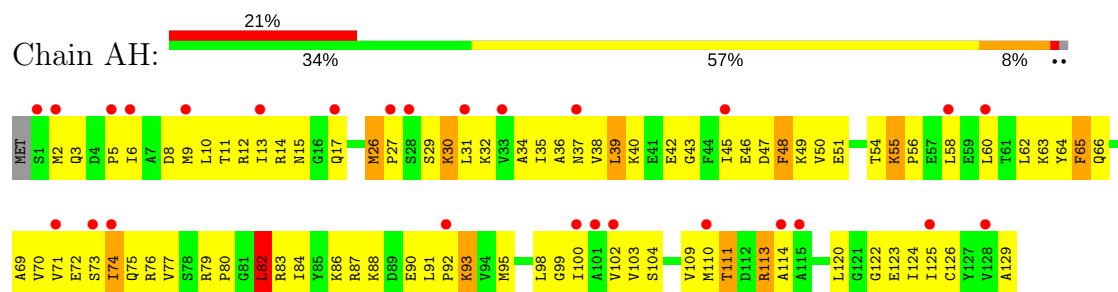




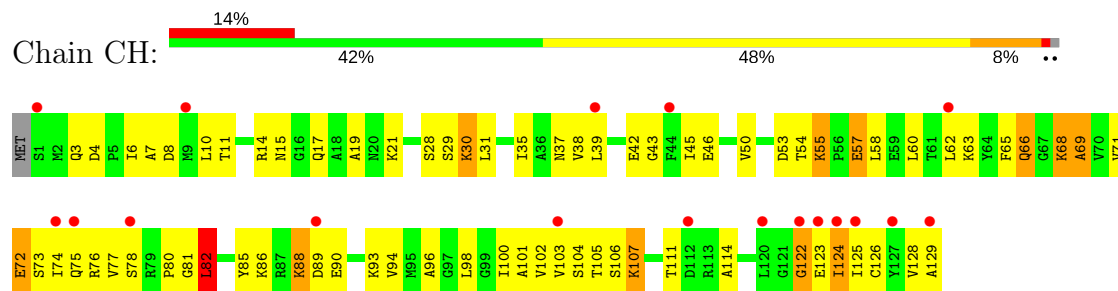
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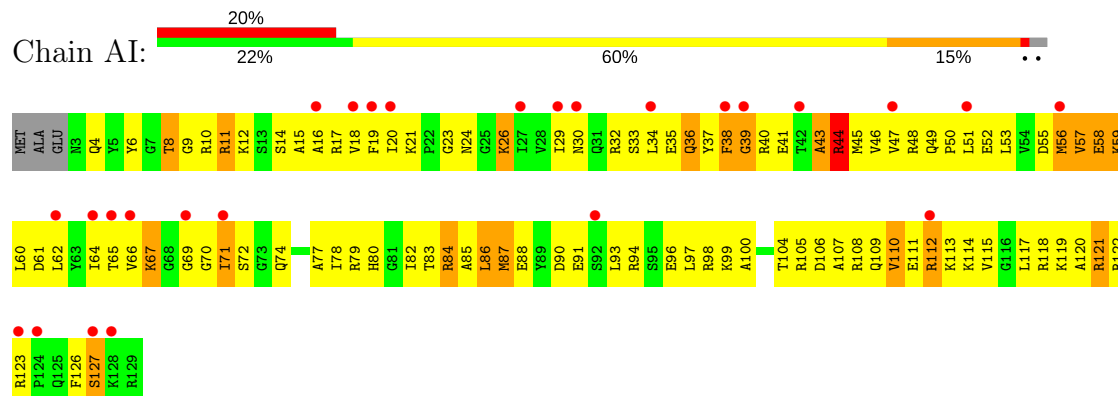
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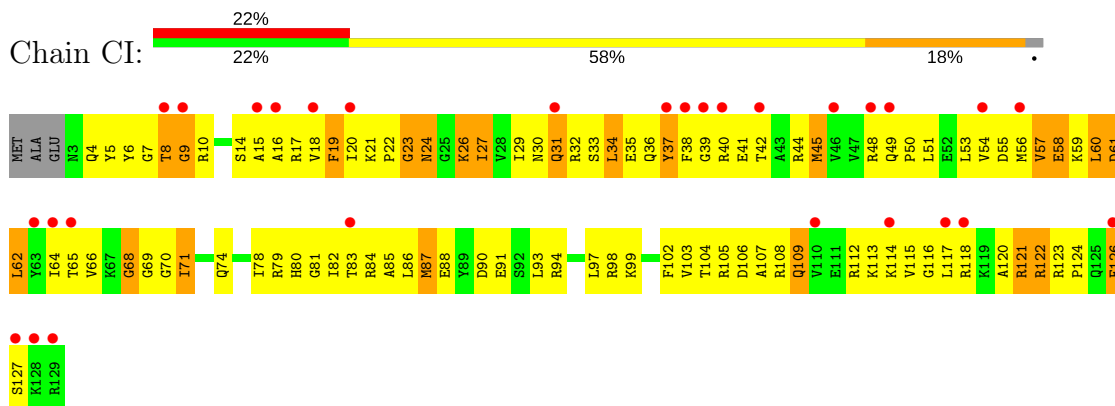
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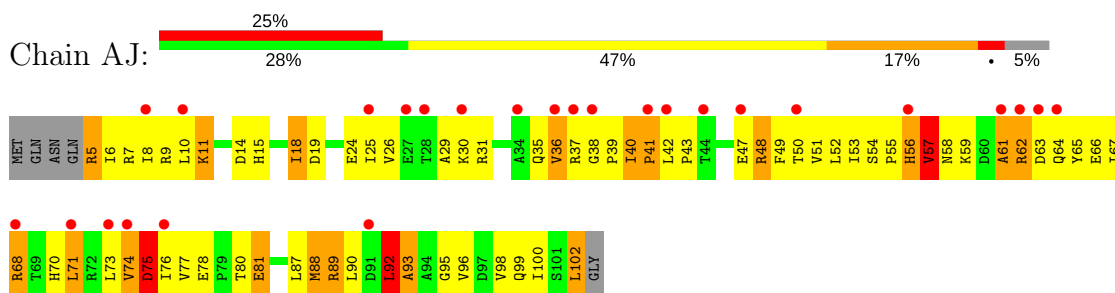
• Molecule 8: 30S RIBOSOMAL PROTEIN S9



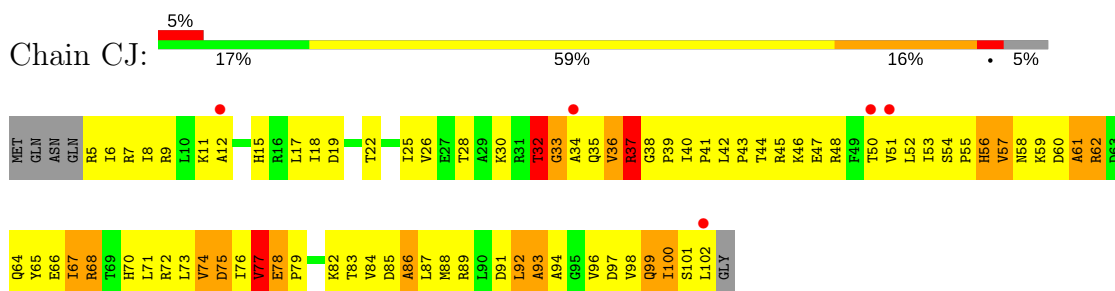
• Molecule 8: 30S RIBOSOMAL PROTEIN S9



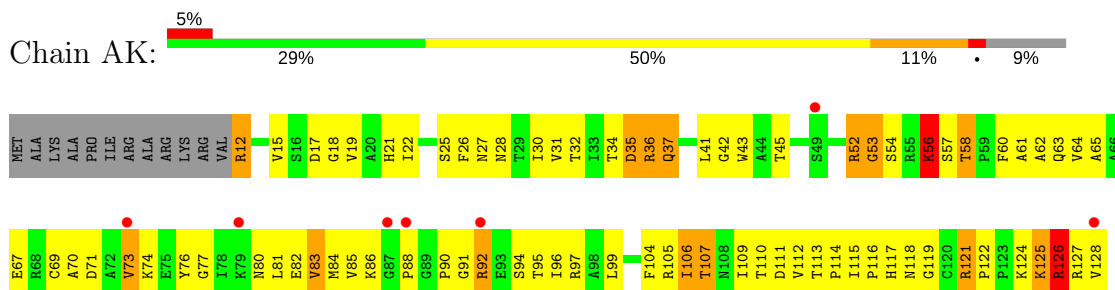
• Molecule 9: 30S RIBOSOMAL PROTEIN S10



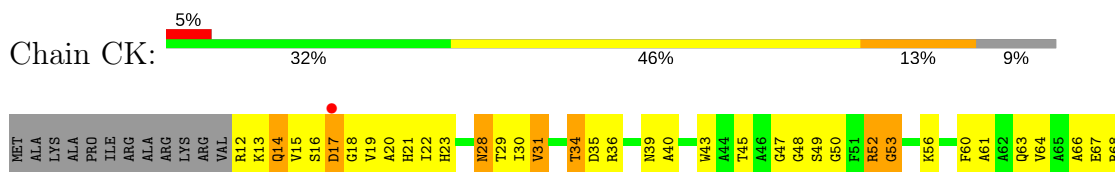
• Molecule 9: 30S RIBOSOMAL PROTEIN S10

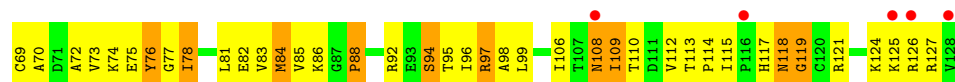


• Molecule 10: 30S RIBOSOMAL PROTEIN S11



• Molecule 10: 30S RIBOSOMAL PROTEIN S11

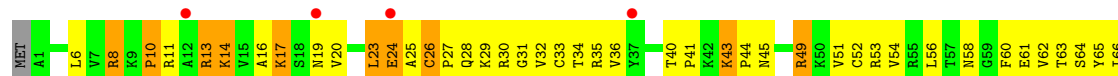




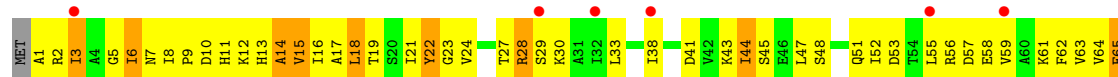
• Molecule 11: 30S RIBOSOMAL PROTEIN S12



• Molecule 11: 30S RIBOSOMAL PROTEIN S12



• Molecule 12: 30S RIBOSOMAL PROTEIN S13

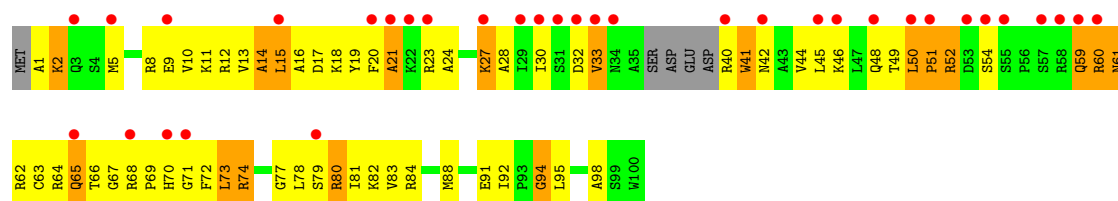


• Molecule 12: 30S RIBOSOMAL PROTEIN S13

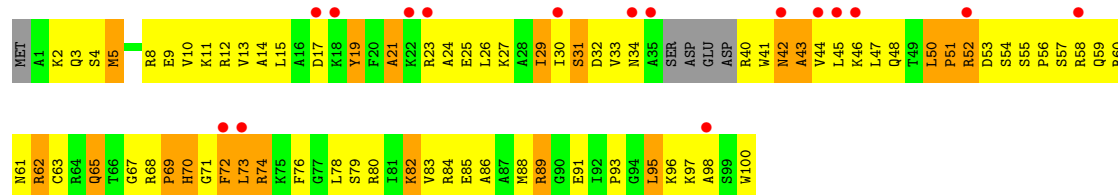


• Molecule 13: 30S RIBOSOMAL PROTEIN S14

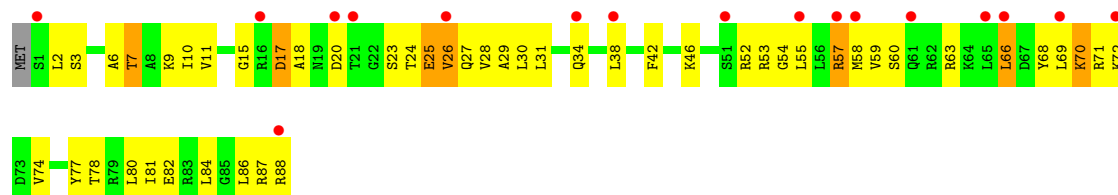




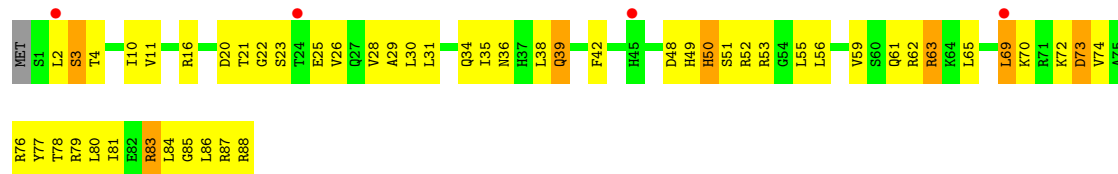
• Molecule 13: 30S RIBOSOMAL PROTEIN S14



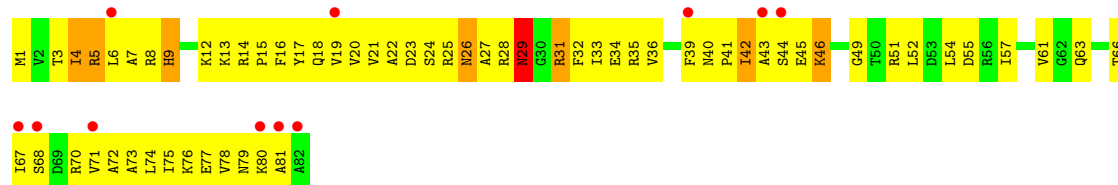
• Molecule 14: 30S RIBOSOMAL PROTEIN S15



• Molecule 14: 30S RIBOSOMAL PROTEIN S15

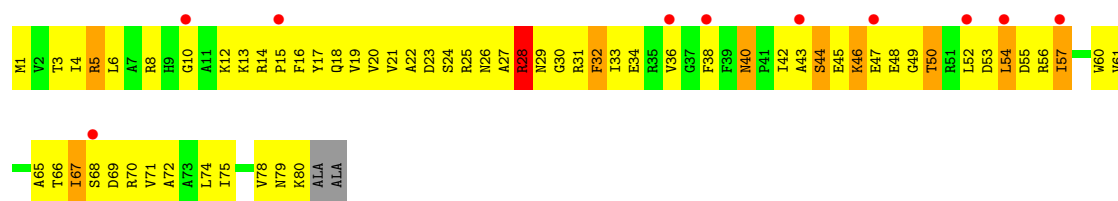


• Molecule 15: 30S RIBOSOMAL PROTEIN S16

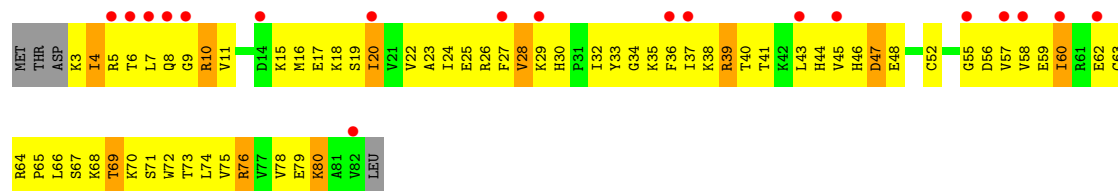


• Molecule 15: 30S RIBOSOMAL PROTEIN S16

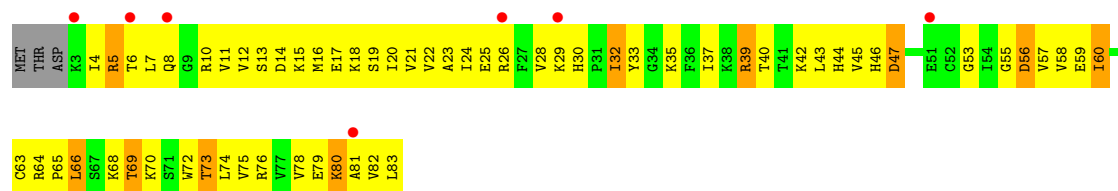




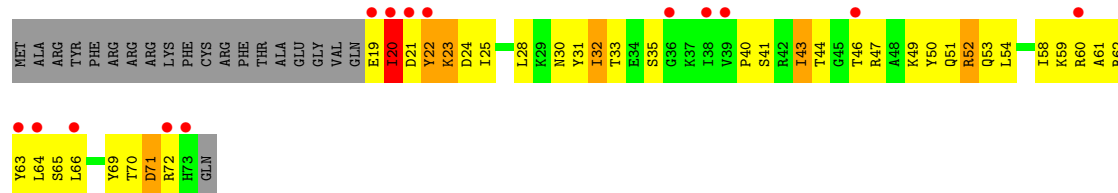
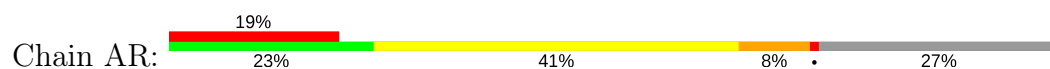
• Molecule 16: 30S RIBOSOMAL PROTEIN S17



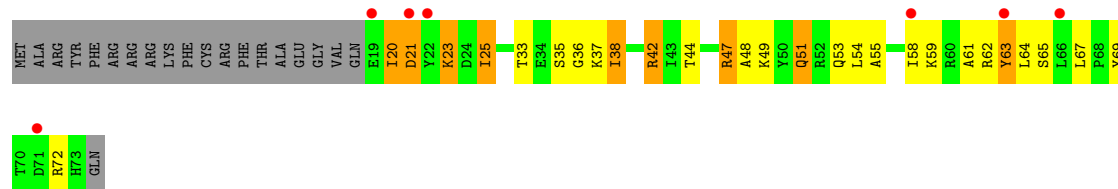
• Molecule 16: 30S RIBOSOMAL PROTEIN S17



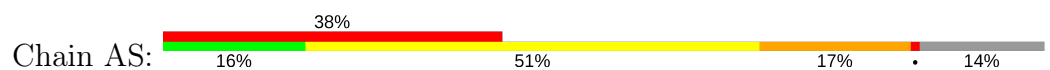
• Molecule 17: 30S RIBOSOMAL PROTEIN S18

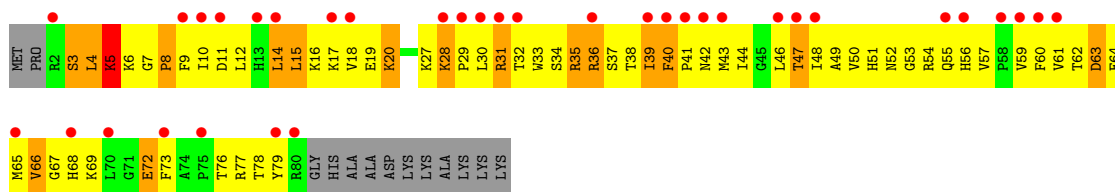


• Molecule 17: 30S RIBOSOMAL PROTEIN S18

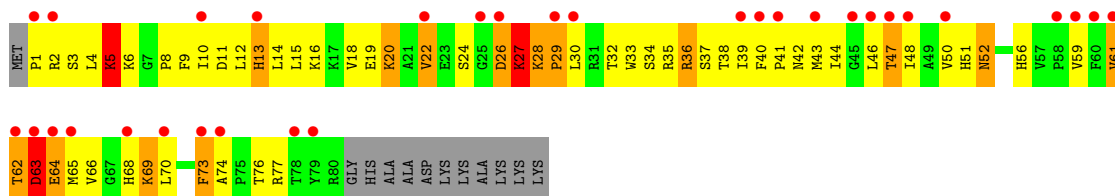


• Molecule 18: 30S RIBOSOMAL PROTEIN S19

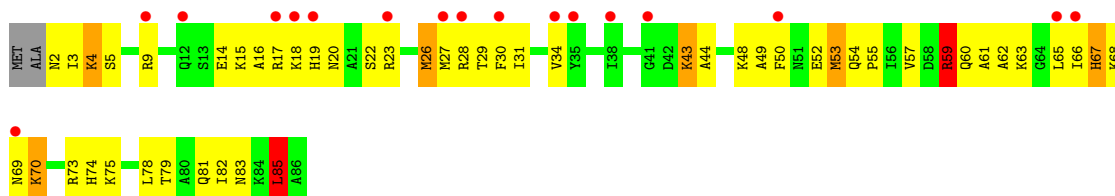




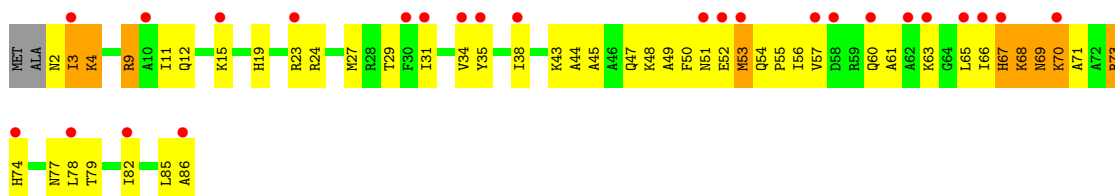
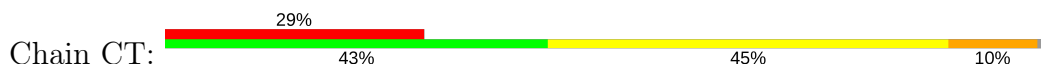
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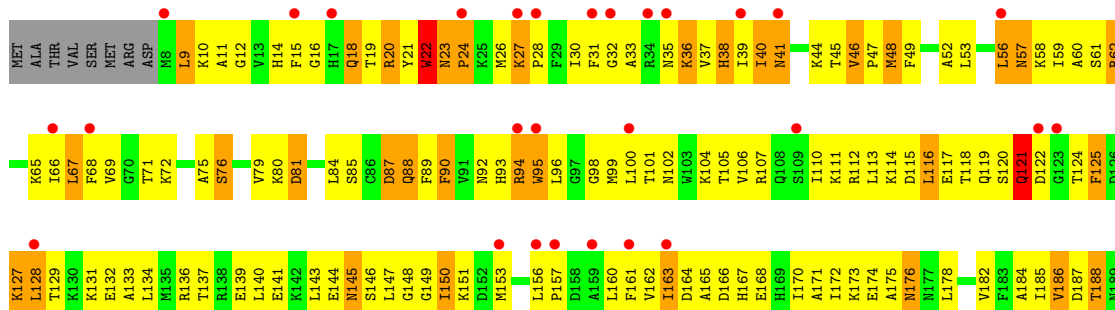
• Molecule 19: 30S RIBOSOMAL PROTEIN S20



• Molecule 19: 30S RIBOSOMAL PROTEIN S20

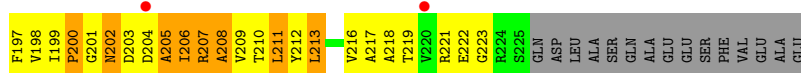
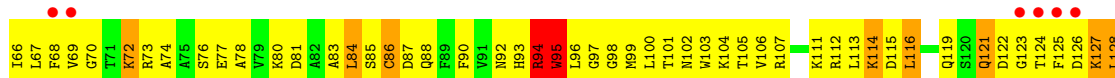
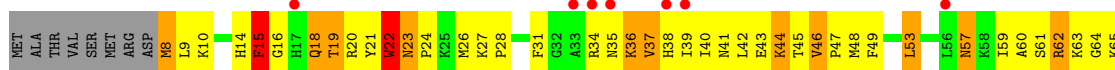


• Molecule 20: 30S RIBOSOMAL PROTEIN S2

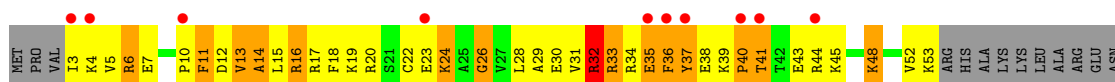




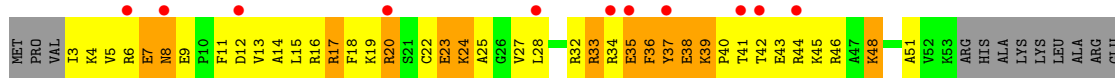
• Molecule 20: 30S RIBOSOMAL PROTEIN S2



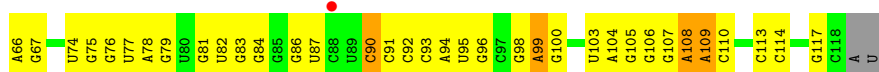
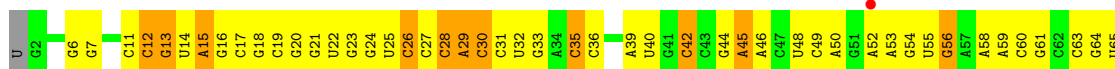
• Molecule 21: 30S RIBOSOMAL PROTEIN S21



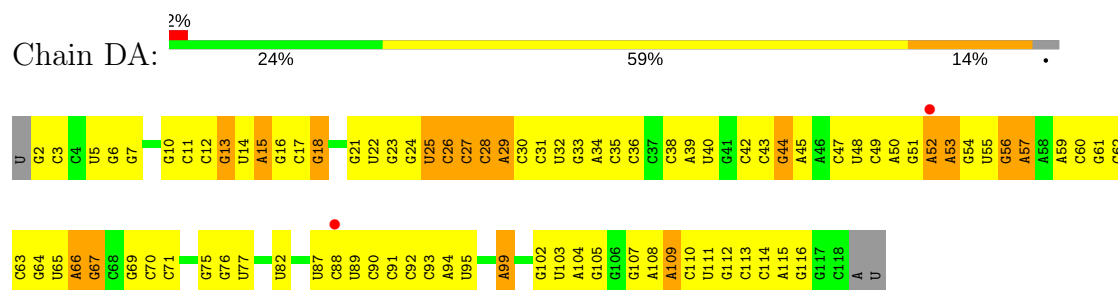
• Molecule 21: 30S RIBOSOMAL PROTEIN S21



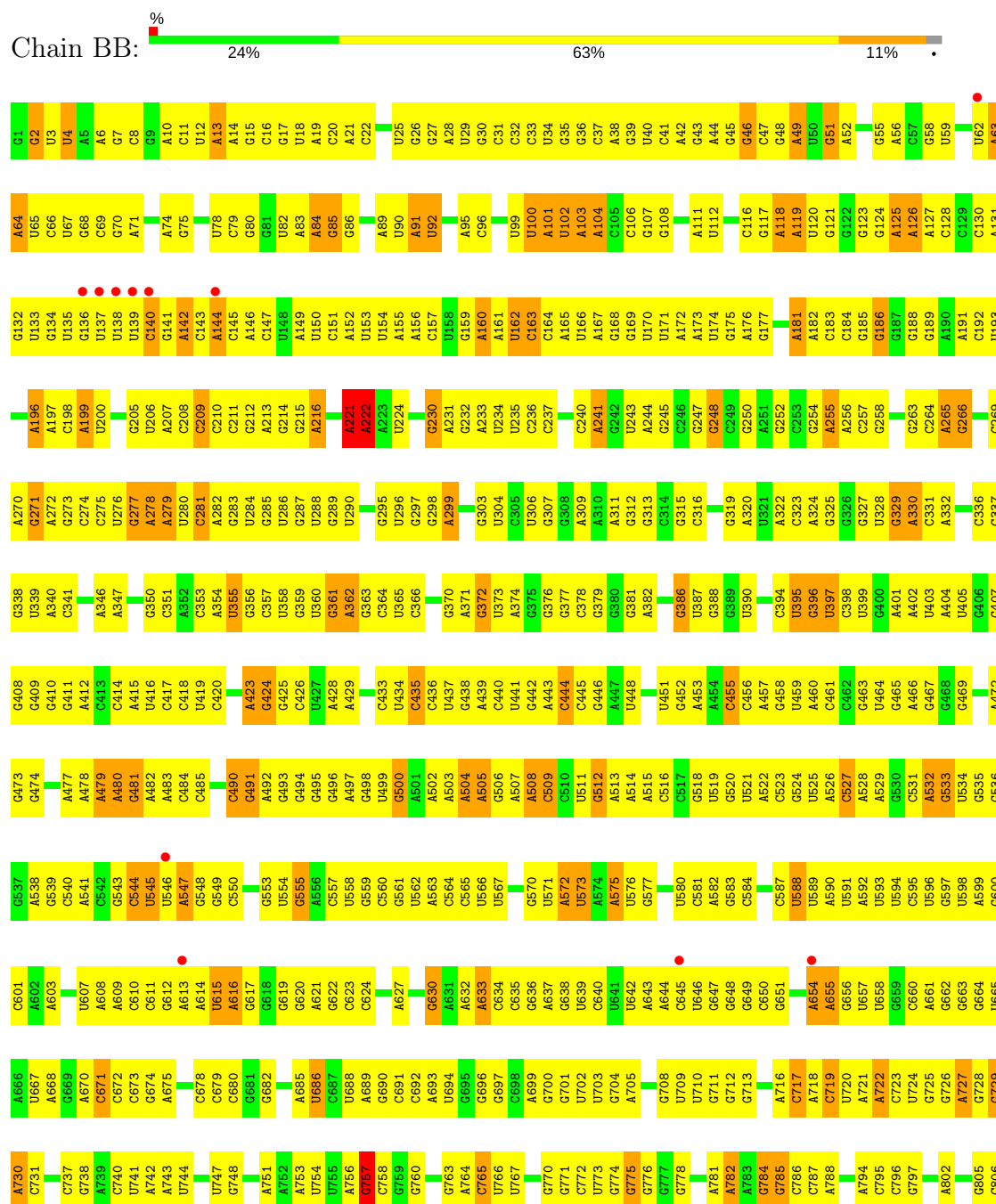
• Molecule 22: 5S RIBOSOMAL RNA



● Molecule 22: 5S RIBOSOMAL RNA

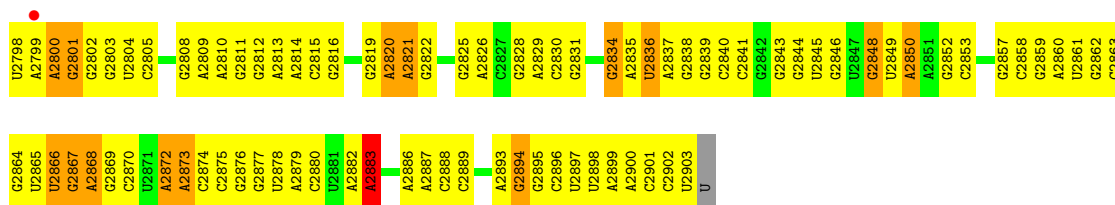


● Molecule 23: 23S RIBOSOMAL RNA

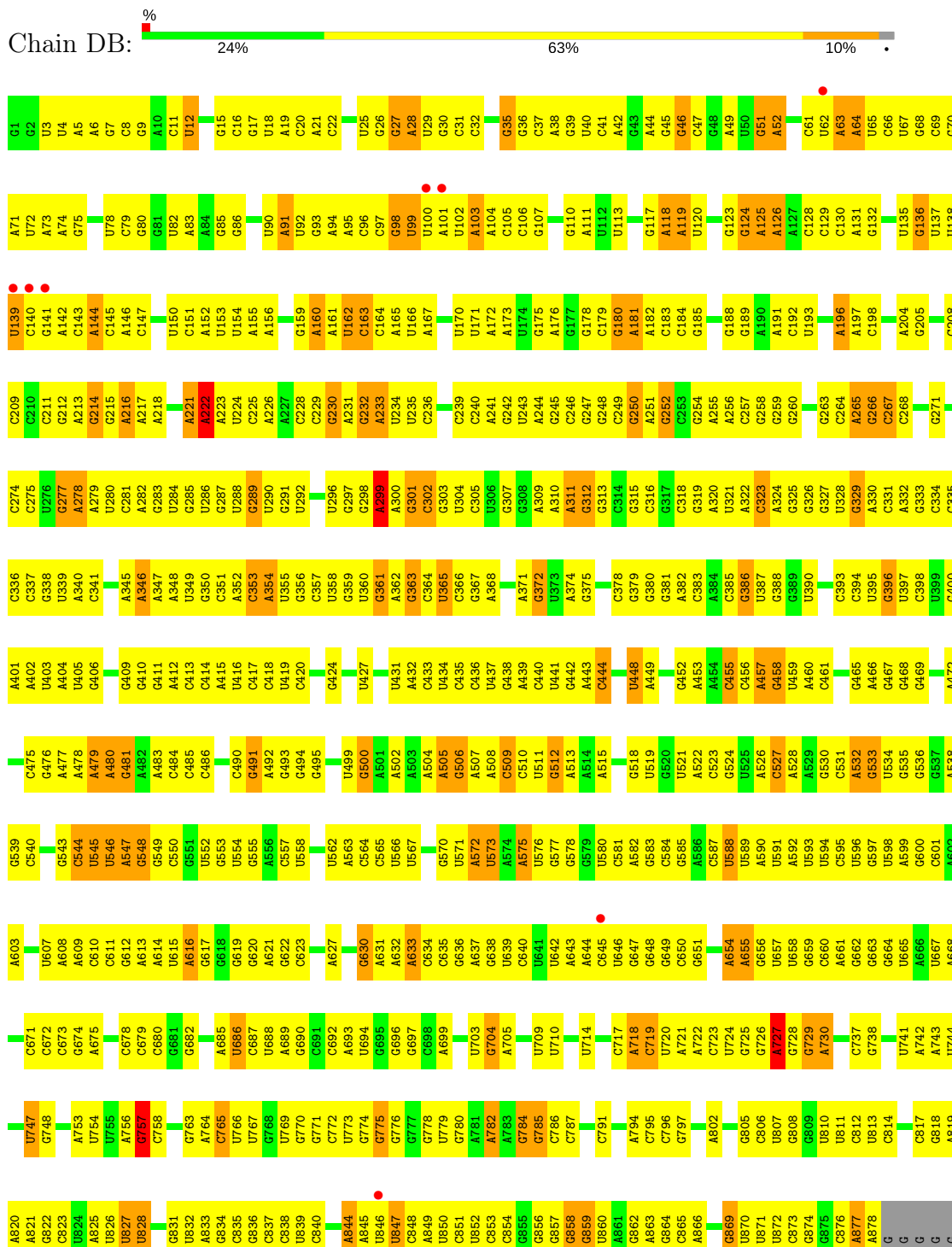


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A1736	U1670	U1594	G1529	C1462	U1401	A1327	A1260	G1197	G1136	G1068	C1006	A877	U813
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G1737	A1596	C1531	U1329	U1464	A1403	U1329	U1263	U1199	G1138	G1070	A943	G	
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G1674	A1597	A1533	U1466	U1466	U1405	C1331	G1266	U1201	A1008	C946	A943	G	
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A1744	C1604	U1539	U1474	U1474	U1411	G1341	G1271	C1207	C1076	A949	A943	G	G822
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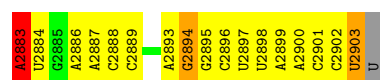


● Molecule 23: 23S RIBOSOMAL RNA

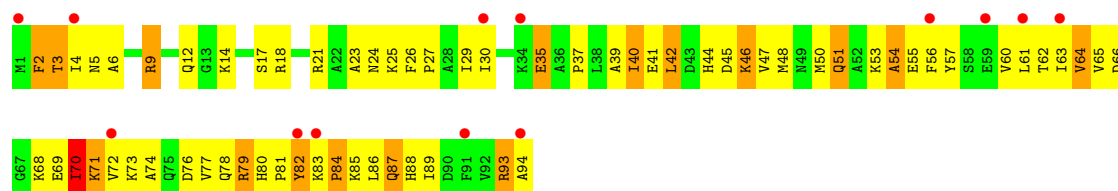




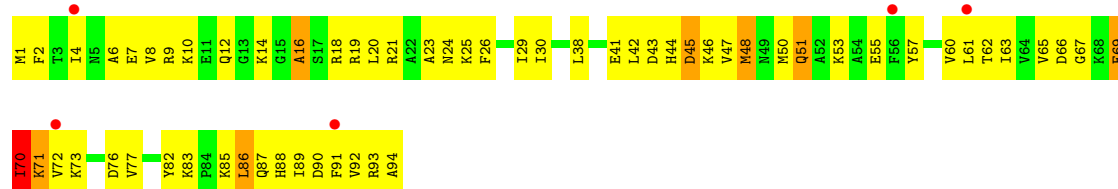
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C2870	C2805	A2736	G2674	G2606	G2535	C2466	U2397	U2332	G2267	U	G2086	G2024	U1951
U2871	U2806	G2737	A2675	U2607	G2536	C2467	U2398	U2333	U2272	U	G2087	C2025	A1952
A2872	G2807	U2738	C2676	C2610	C2537	A2468	G2400	U2334	A2273	U	A2088	U2026	A1953
A2873	U2808	U2739	G2677	G2611	C2538	G2471	U2401	A2335	A2274	U	C2089	G2027	G1954
C2874	A2809	A2740	C2678	U2613	C2540	G2472	U2402	U2336	C2275	U	A2090	U2028	U1955
C2875	G2810	U2741	U2679	U2614	A2541	C2473	C2403	G2337	G2276	U	C2091	G2029	U1956
G2876	G2811	G2742	U2680	C2615	G2542	C2474	U2404	C2338	G2277	U	U2092	G2030	U1957
C2877	A2812	U2743	C2681	G2616	G2543	A2475	G2405	A2339	A2278	U	A2093	A2031	C1958
U2878	U2813	G2744	A2682	U2617	G2544	A2476	A2406	G2341	G2279	U	A2094	G2032	G1959
A2879	C2814	C2745	G2683	G2618	U2545	A2477	A2407	C2342	A2280	U	A2095	A2033	G1964
G2880	U2815	U2746	U2684	U2619	U2546	A2478	U2408	U2343	A2281	U	C2096	G2034	
U2881	G2816	G2747	G2685	G2621	A2547	U2479	U2409	G2344	G2282	U	U2097	U2035	
A2882	U2817	U2748	G2686	C2626	U2548	C2480	G2410	G2345	C2283	U	A2098		
	G2819	G2751	U2687	C2626		G2481							



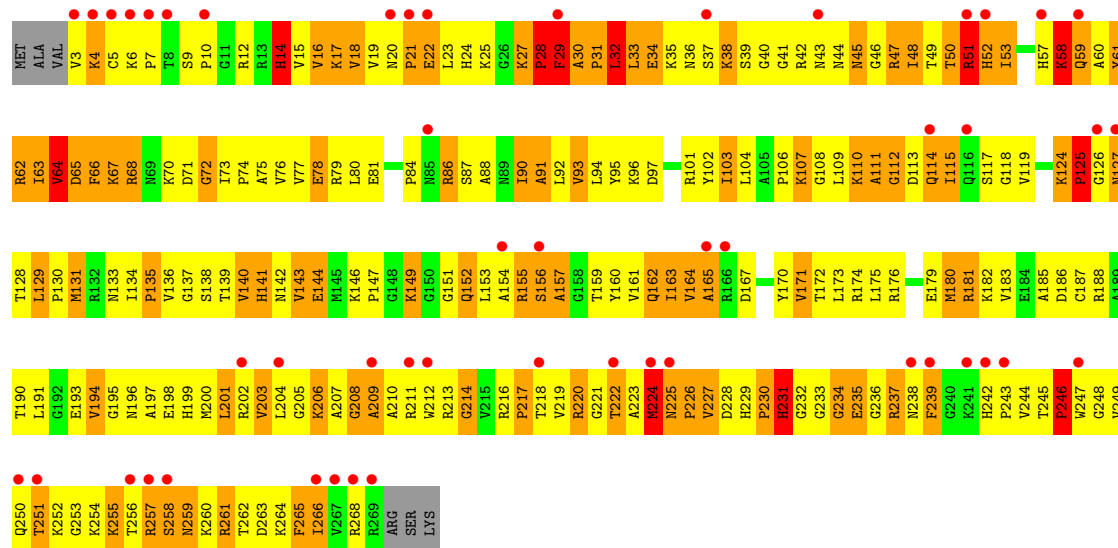
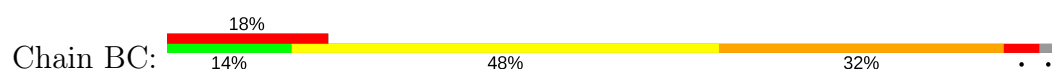
• Molecule 24: 50S RIBOSOMAL PROTEIN L25



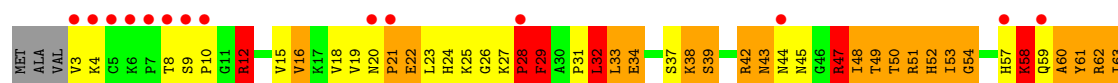
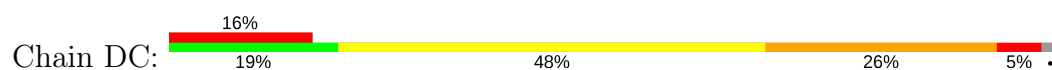
• Molecule 24: 50S RIBOSOMAL PROTEIN L25

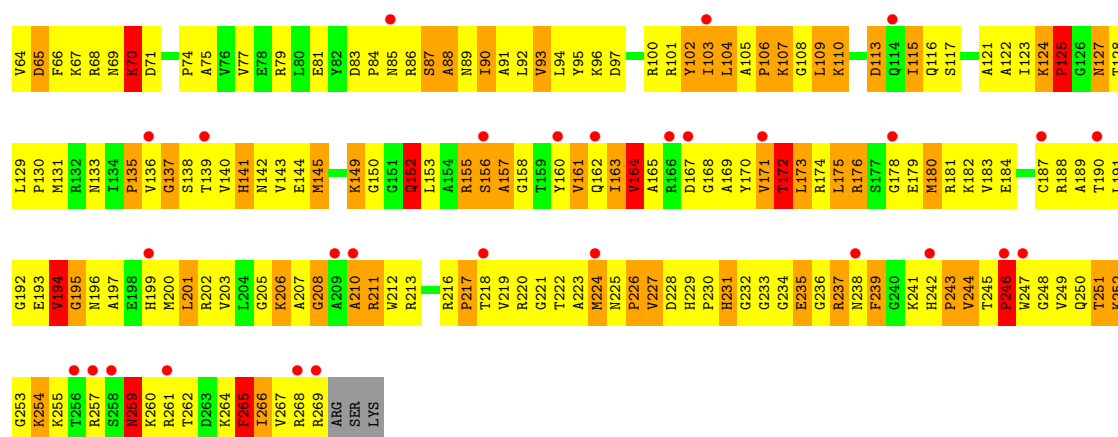


• Molecule 25: 50S RIBOSOMAL PROTEIN L2

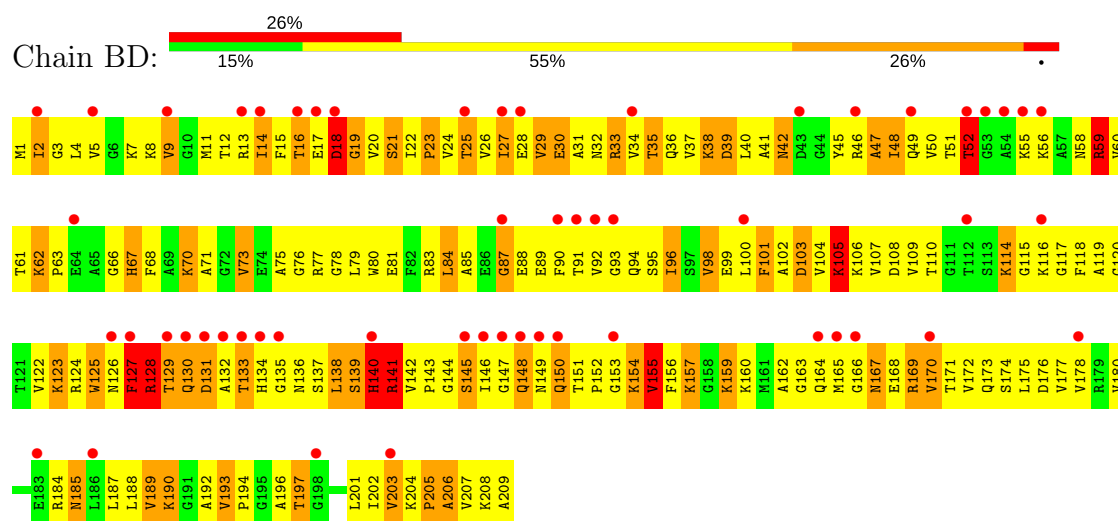


• Molecule 25: 50S RIBOSOMAL PROTEIN L2

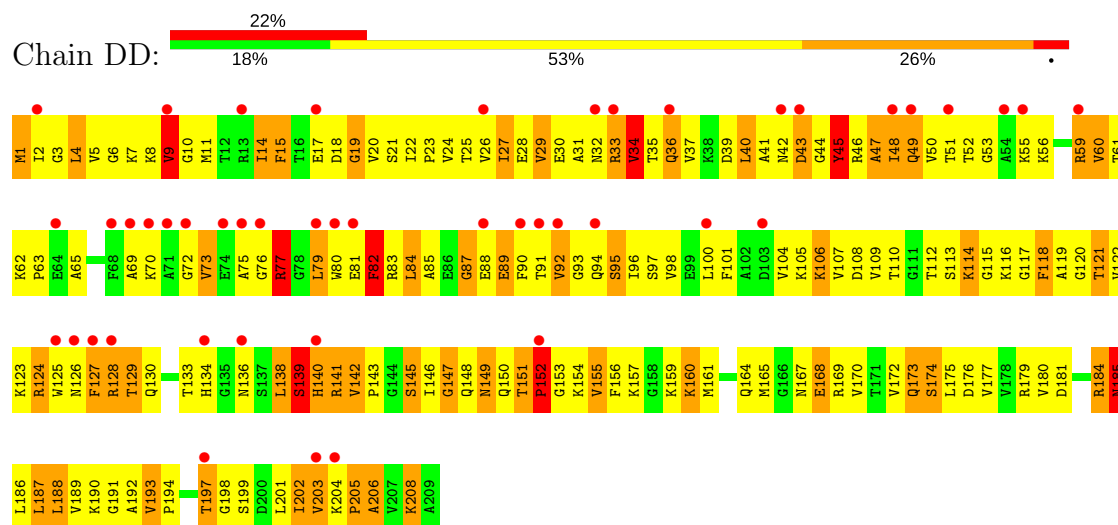




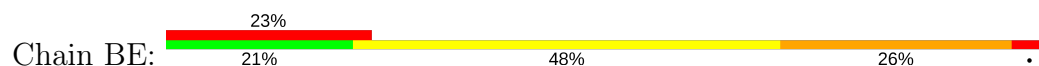
● Molecule 26: 50S RIBOSOMAL PROTEIN L3

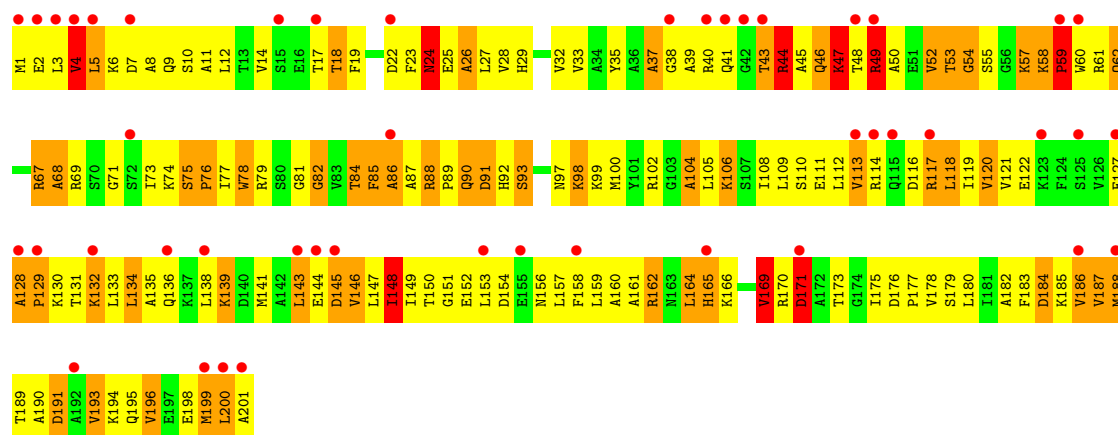


● Molecule 26: 50S RIBOSOMAL PROTEIN L3



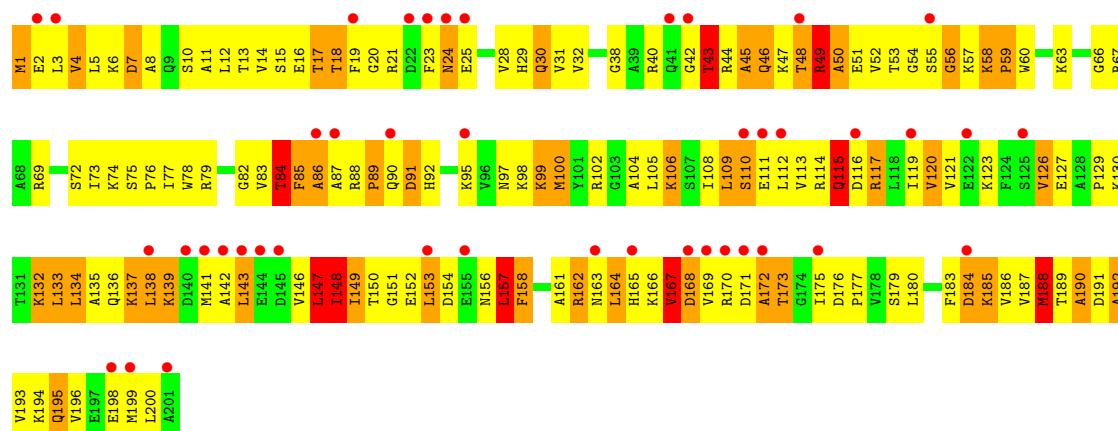
● Molecule 27: 50S RIBOSOMAL PROTEIN L4





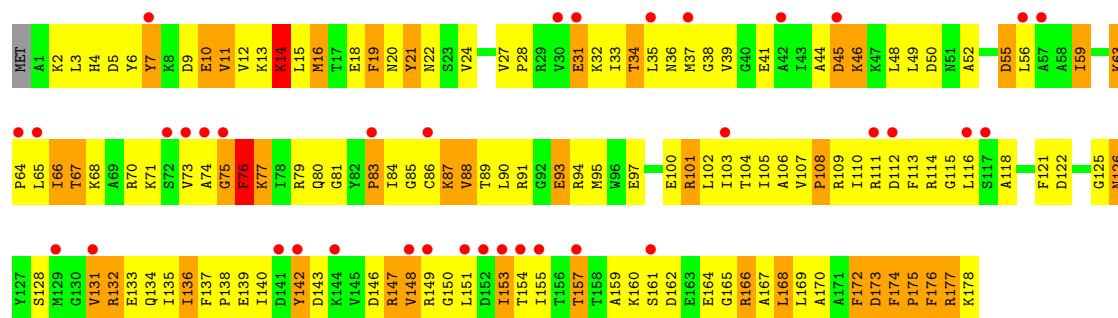
• Molecule 27: 50S RIBOSOMAL PROTEIN L4

Chain DE: 21% 22% 51% 23%



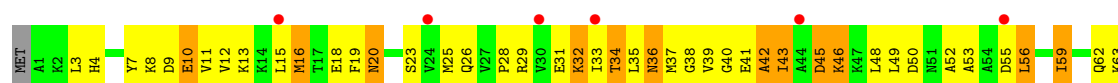
• Molecule 28: 50S RIBOSOMAL PROTEIN L5

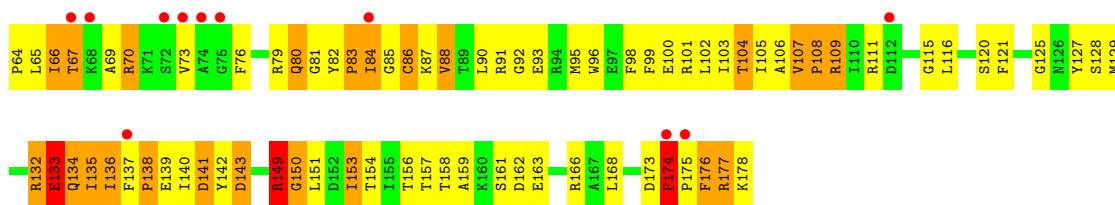
Chain BF: 20% 25% 51% 22%



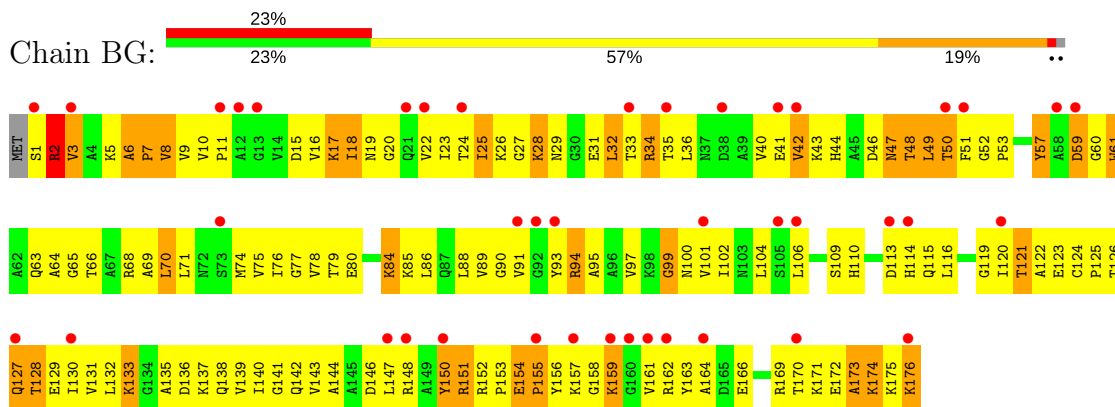
• Molecule 28: 50S RIBOSOMAL PROTEIN L5

Chain DF: 9% 32% 46% 20%

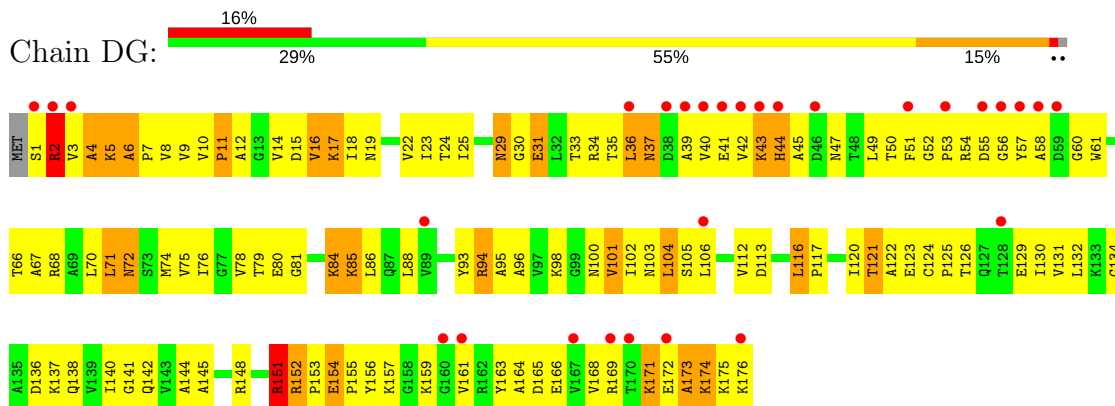




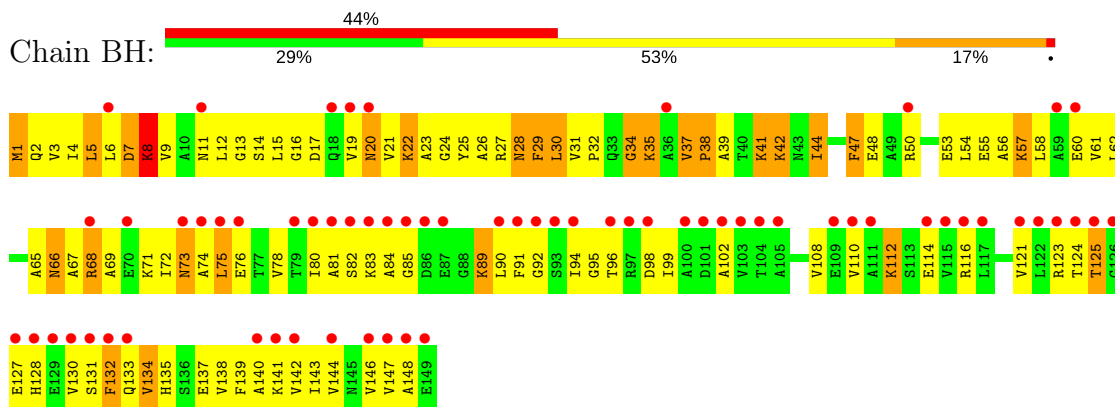
• Molecule 29: 50S RIBOSOMAL PROTEIN L6



• Molecule 29: 50S RIBOSOMAL PROTEIN L6

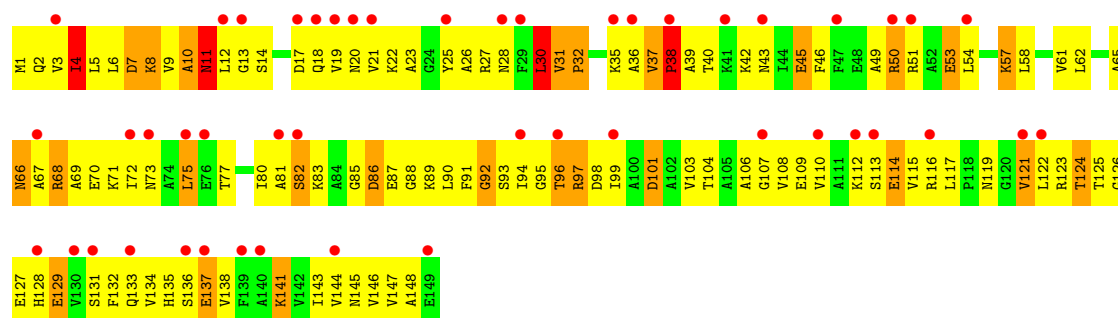


• Molecule 30: 50S RIBOSOMAL PROTEIN L9

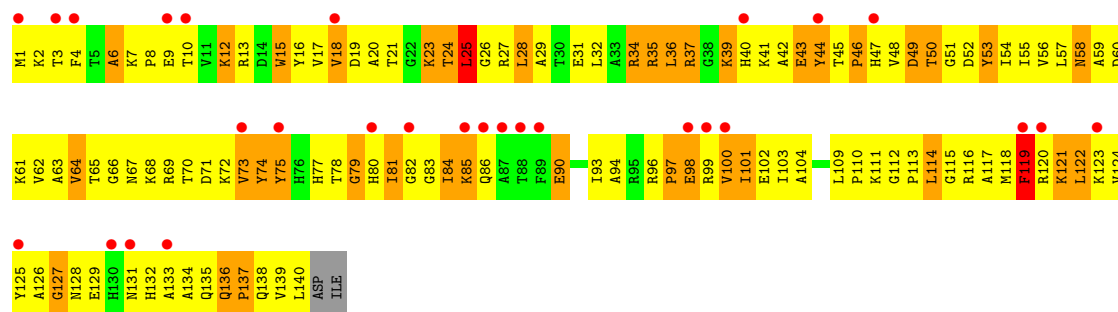
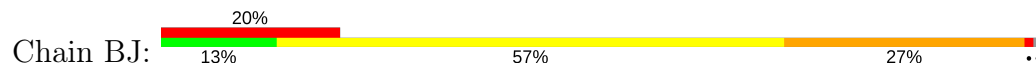


• Molecule 30: 50S RIBOSOMAL PROTEIN L9

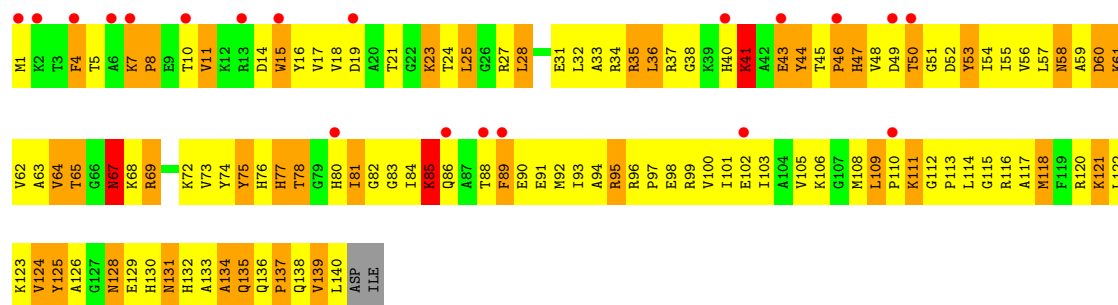
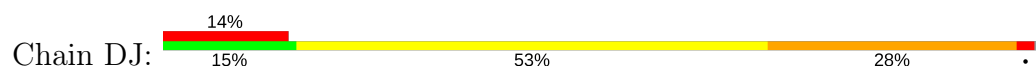




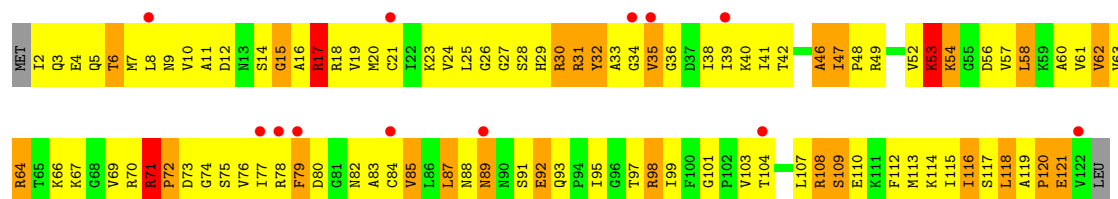
• Molecule 31: 50S RIBOSOMAL PROTEIN L13



• Molecule 31: 50S RIBOSOMAL PROTEIN L13

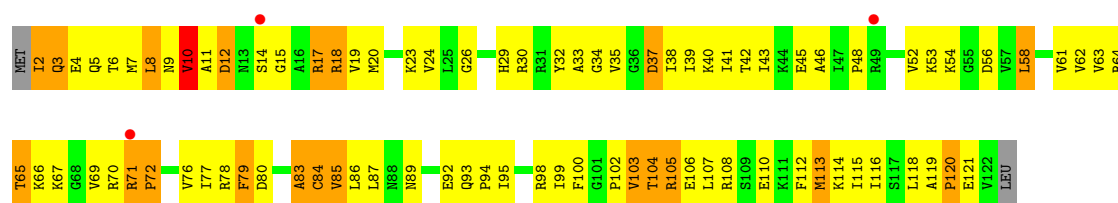


• Molecule 32: 50S RIBOSOMAL PROTEIN L14

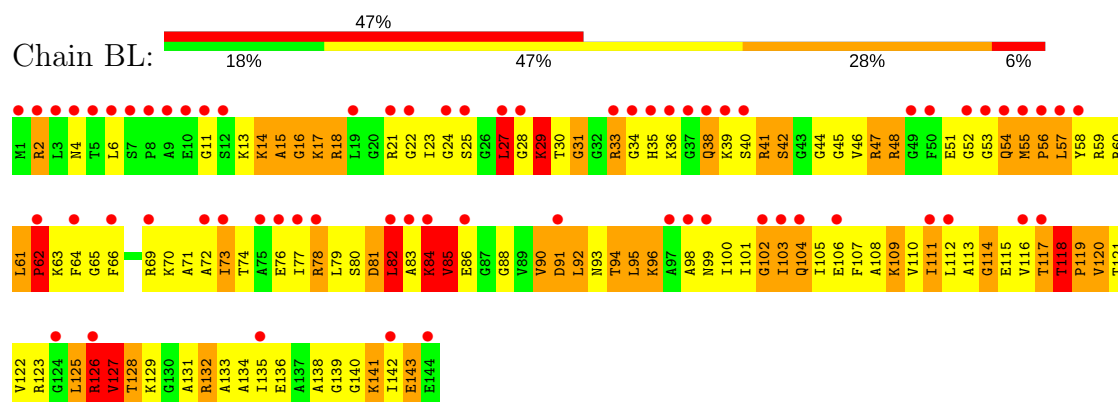


• Molecule 32: 50S RIBOSOMAL PROTEIN L14

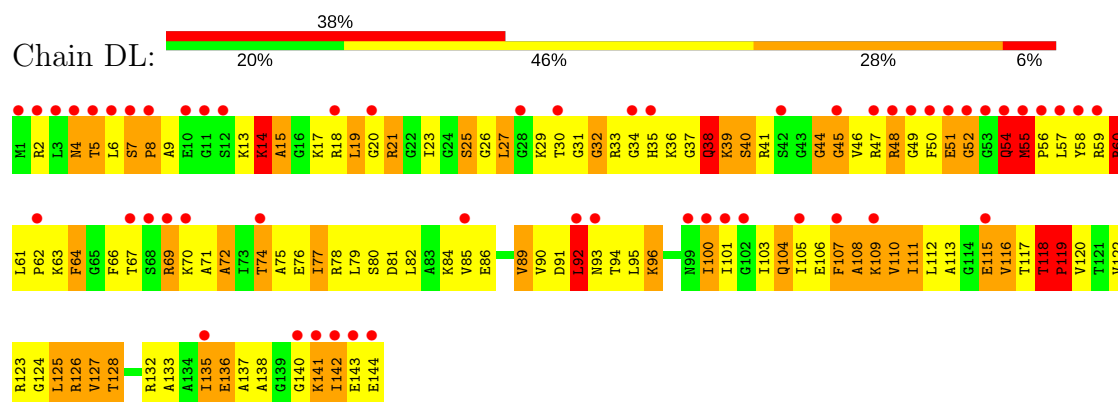




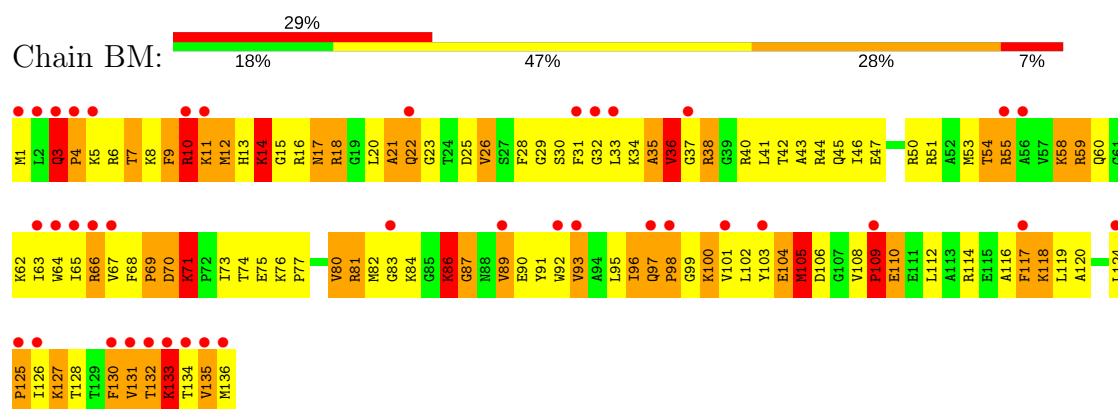
• Molecule 33: 50S RIBOSOMAL PROTEIN L15



• Molecule 33: 50S RIBOSOMAL PROTEIN L15

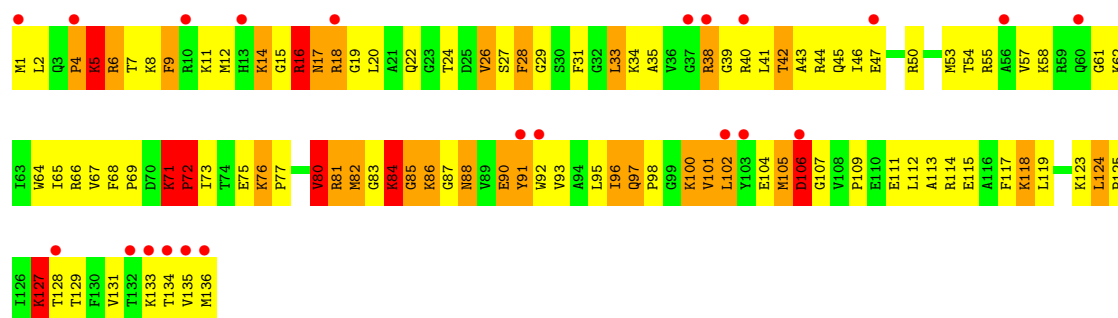


• Molecule 34: 50S RIBOSOMAL PROTEIN L16

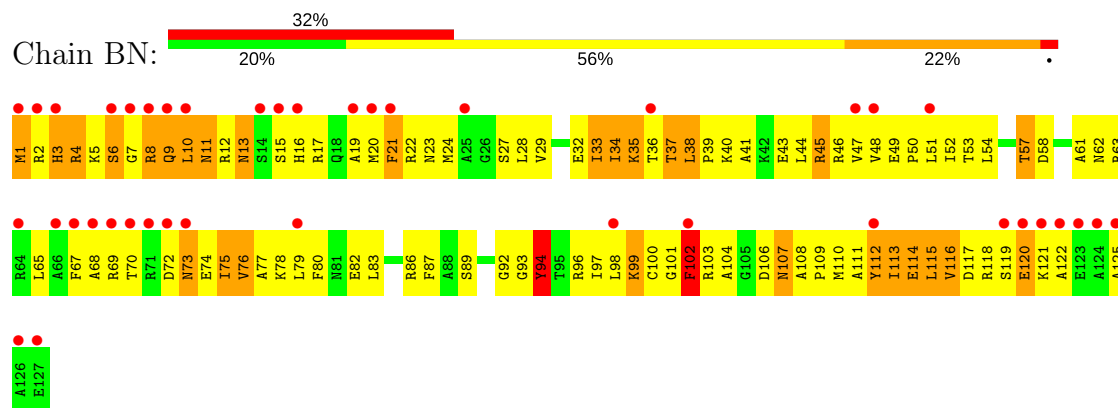


• Molecule 34: 50S RIBOSOMAL PROTEIN L16

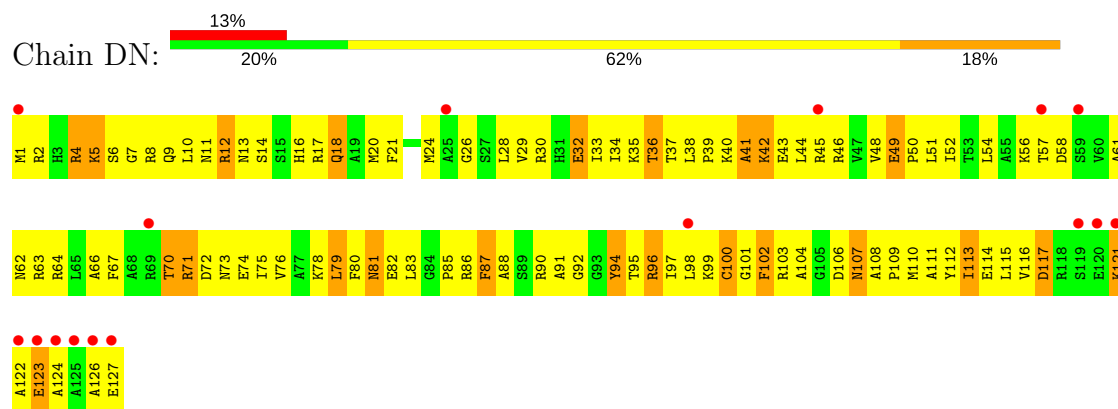




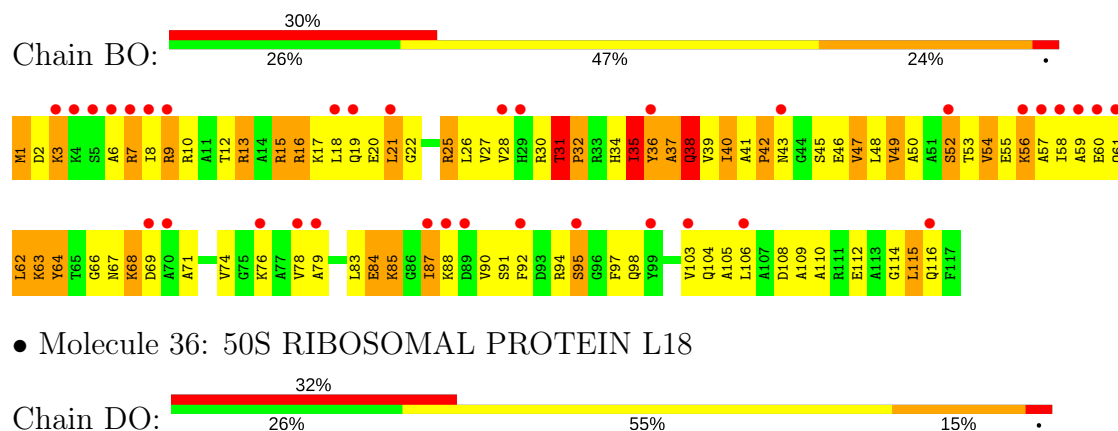
• Molecule 35: 50S RIBOSOMAL PROTEIN L17



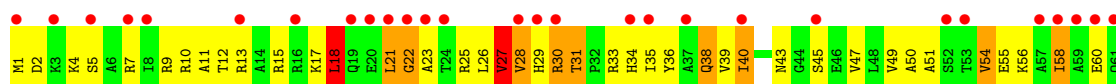
• Molecule 35: 50S RIBOSOMAL PROTEIN L17



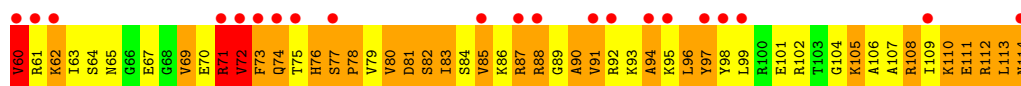
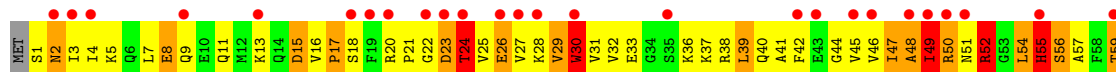
• Molecule 36: 50S RIBOSOMAL PROTEIN L18



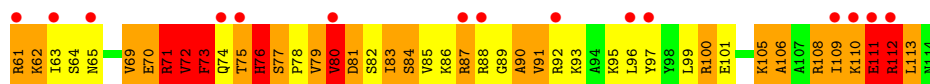
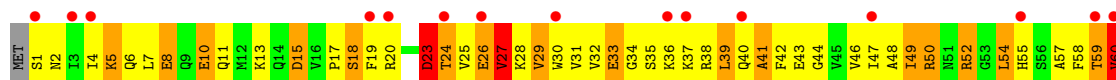
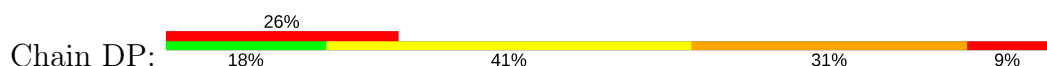
• Molecule 36: 50S RIBOSOMAL PROTEIN L18



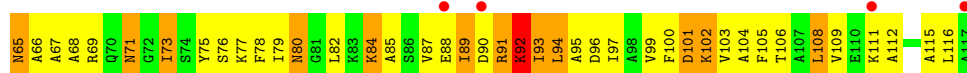
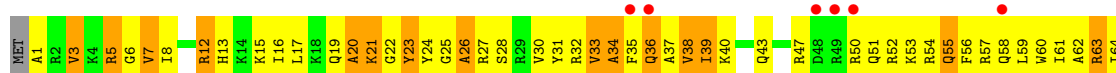
• Molecule 37: 50S RIBOSOMAL PROTEIN L19



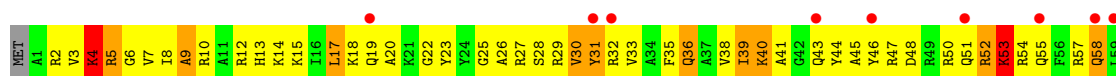
• Molecule 37: 50S RIBOSOMAL PROTEIN L19



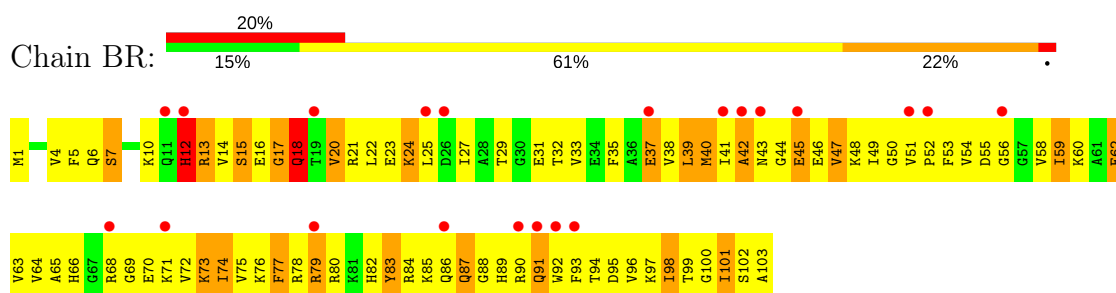
• Molecule 38: 50S RIBOSOMAL PROTEIN L20



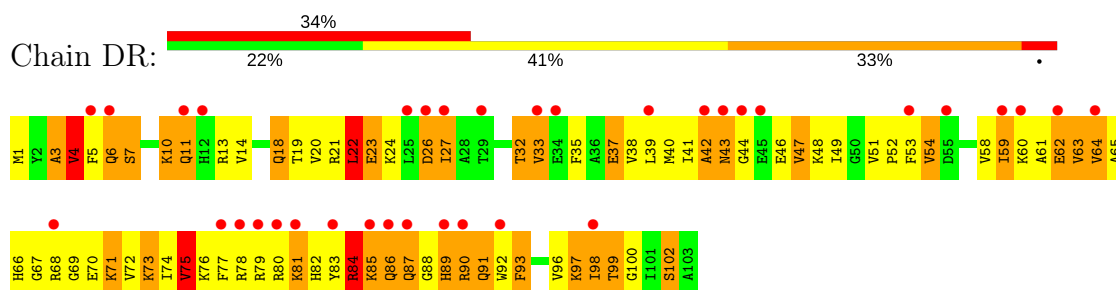
• Molecule 38: 50S RIBOSOMAL PROTEIN L20



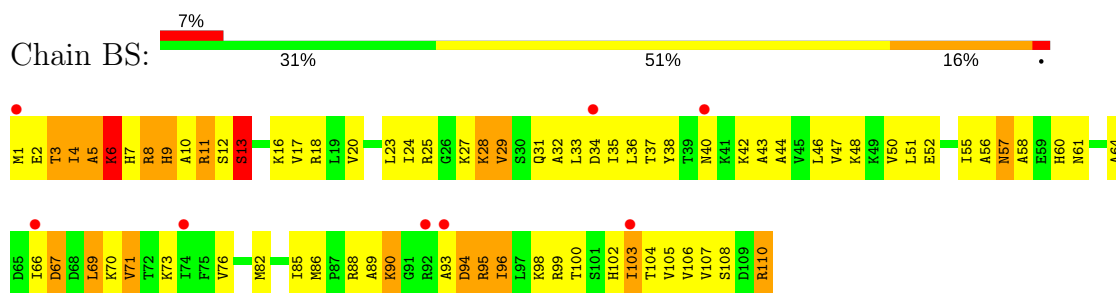
• Molecule 39: 50S RIBOSOMAL PROTEIN L21



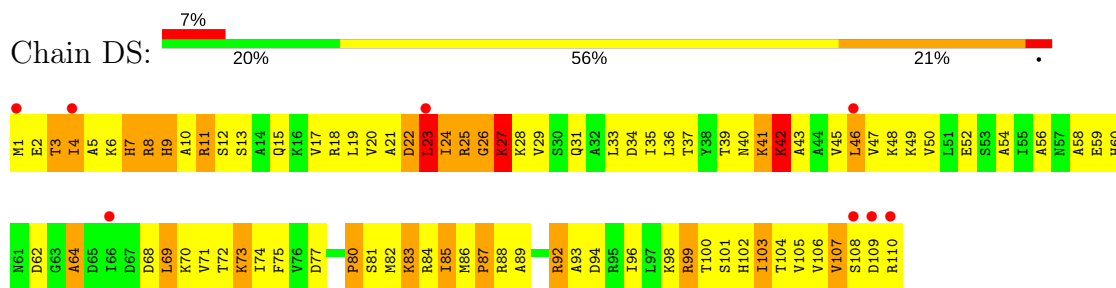
• Molecule 39: 50S RIBOSOMAL PROTEIN L21



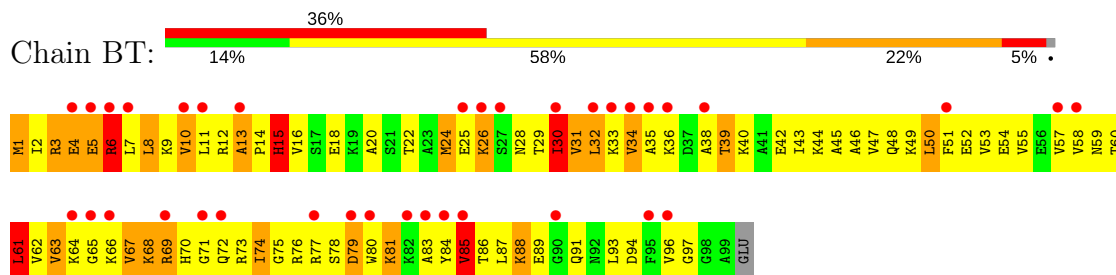
• Molecule 40: 50S RIBOSOMAL PROTEIN L22



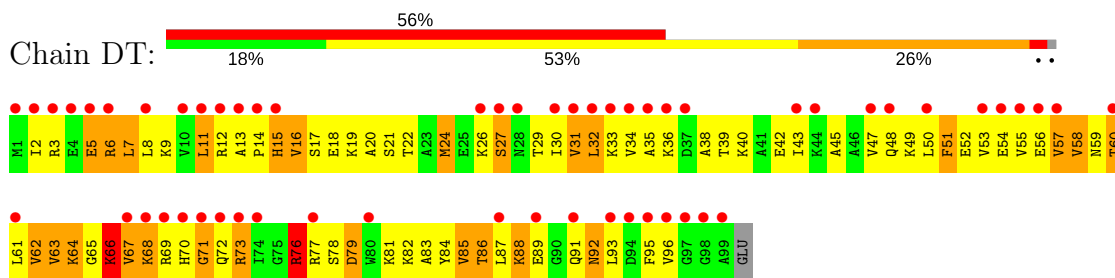
• Molecule 40: 50S RIBOSOMAL PROTEIN L22



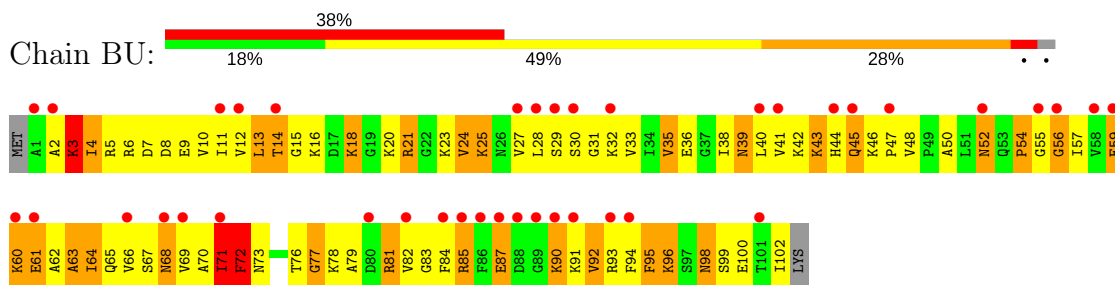
• Molecule 41: 50S RIBOSOMAL PROTEIN L23



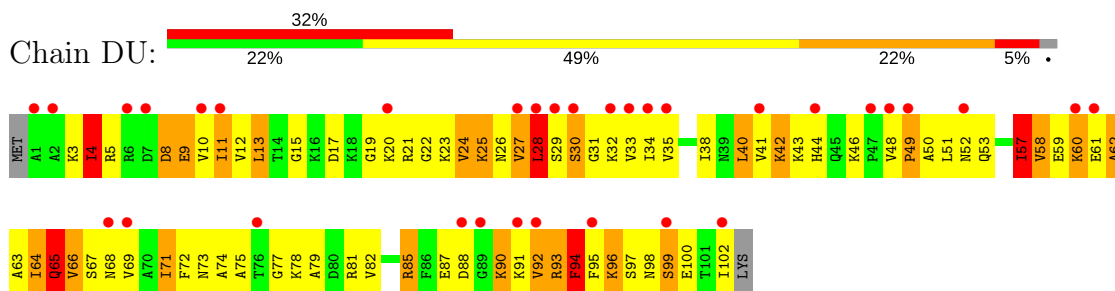
• Molecule 41: 50S RIBOSOMAL PROTEIN L23



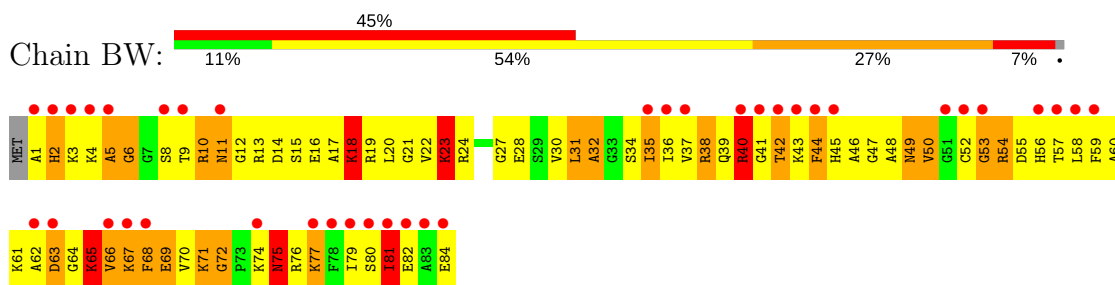
● Molecule 42: 50S RIBOSOMAL PROTEIN L24



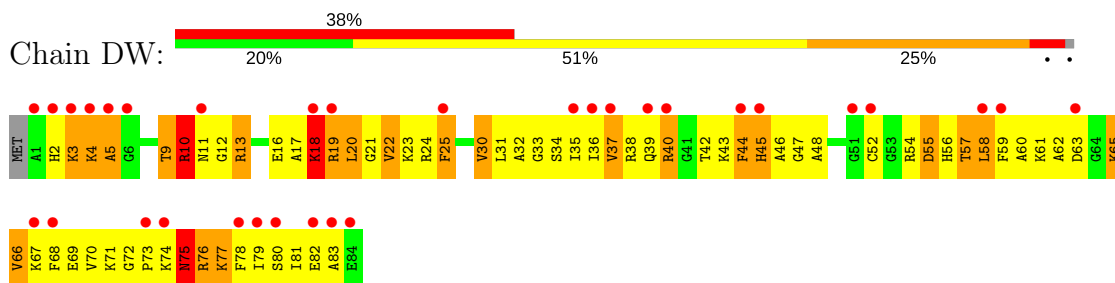
- Molecule 42: 50S RIBOSOMAL PROTEIN L24



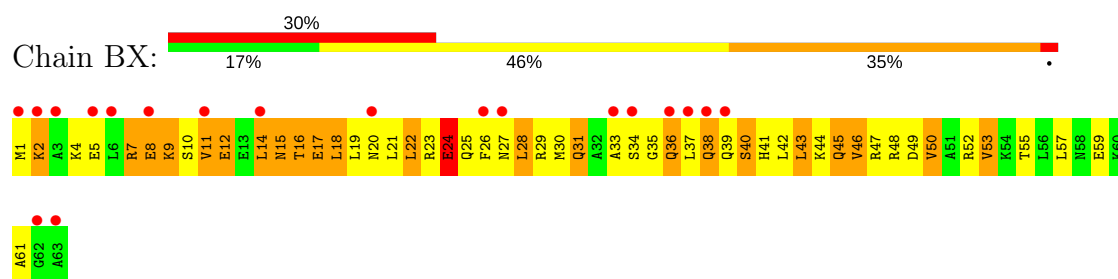
● Molecule 43: 50S RIBOSOMAL PROTEIN L27



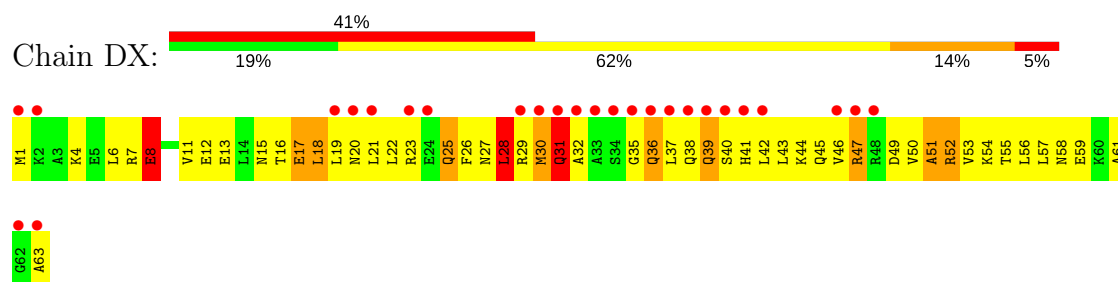
● Molecule 43: 50S RIBOSOMAL PROTEIN L27



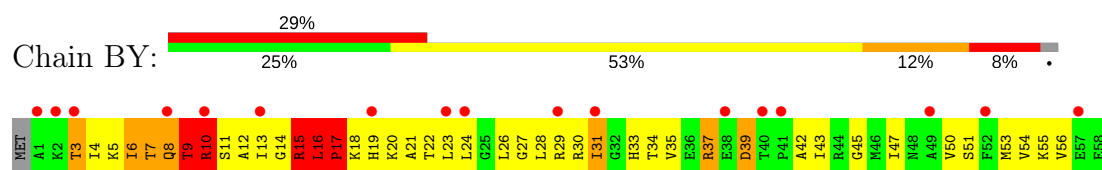
- Molecule 44: 50S RIBOSOMAL PROTEIN L29



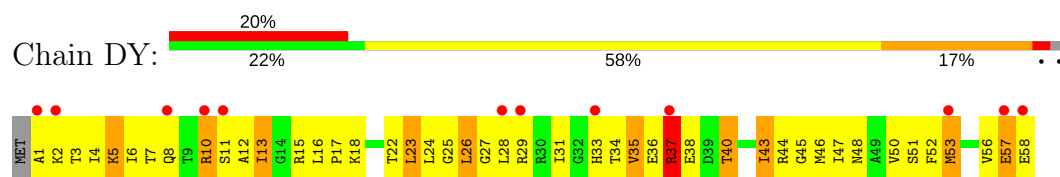
● Molecule 44: 50S RIBOSOMAL PROTEIN L29



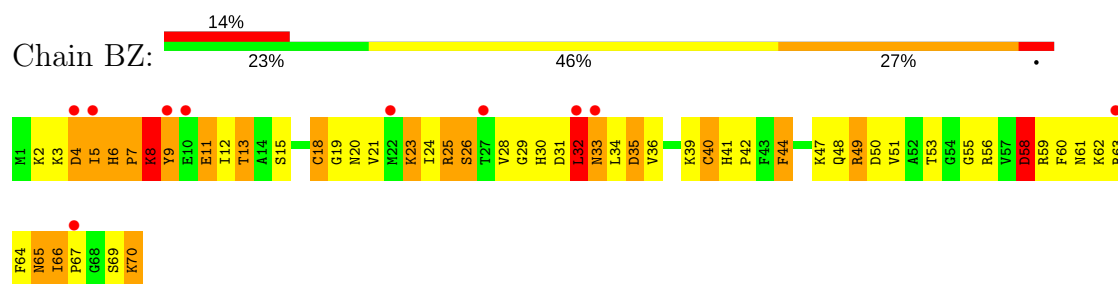
● Molecule 45: 50S RIBOSOMAL PROTEIN L30



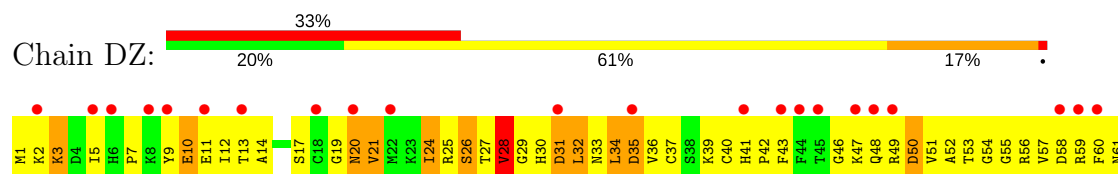
● Molecule 45: 50S RIBOSOMAL PROTEIN L30

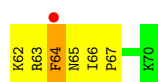


● Molecule 46: 50S RIBOSOMAL PROTEIN L31

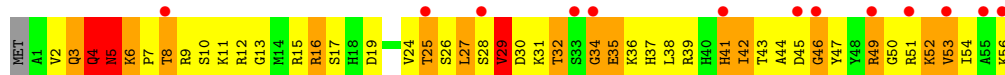
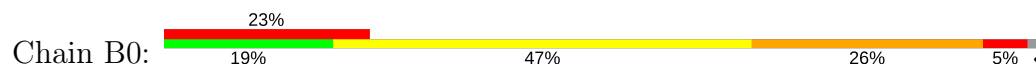


● Molecule 46: 50S RIBOSOMAL PROTEIN L31





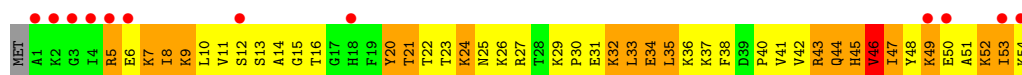
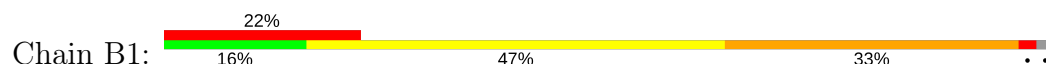
• Molecule 47: 50S RIBOSOMAL PROTEIN L32



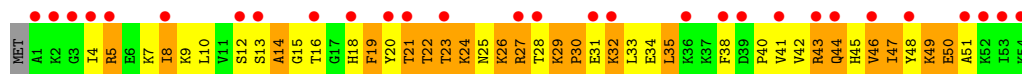
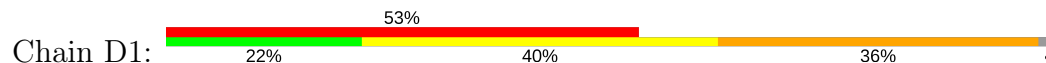
• Molecule 47: 50S RIBOSOMAL PROTEIN L32



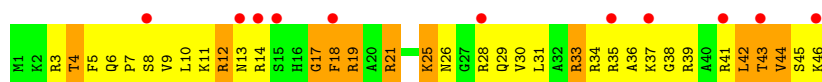
• Molecule 48: 50S RIBOSOMAL PROTEIN L33



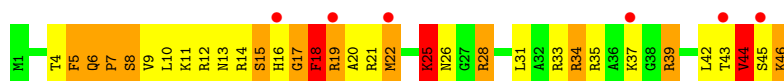
• Molecule 48: 50S RIBOSOMAL PROTEIN L33



• Molecule 49: 50S RIBOSOMAL PROTEIN L34

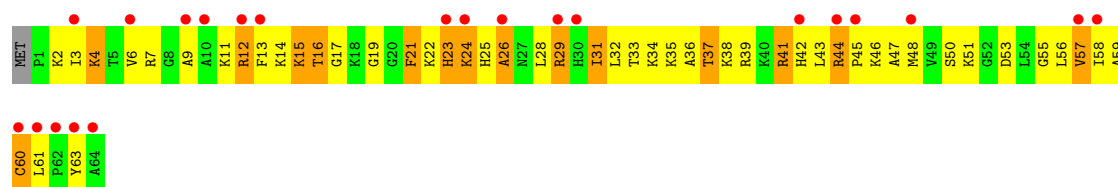


• Molecule 49: 50S RIBOSOMAL PROTEIN L34

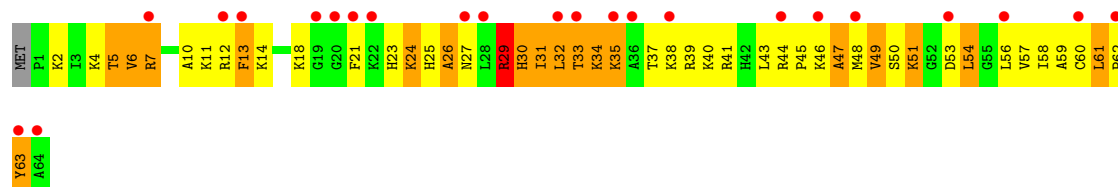


• Molecule 50: 50S RIBOSOMAL PROTEIN L35

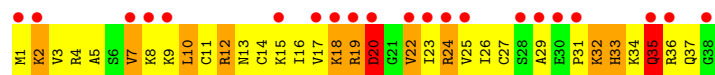
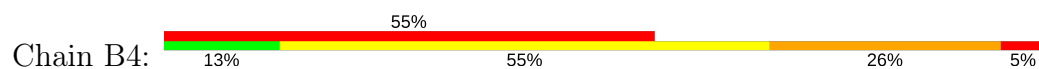




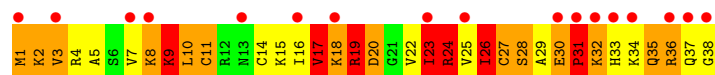
• Molecule 50: 50S RIBOSOMAL PROTEIN L35



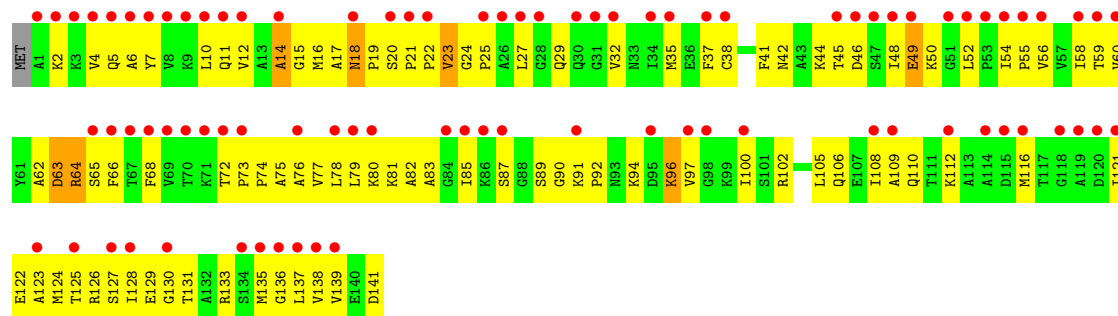
• Molecule 51: 50S RIBOSOMAL PROTEIN L36



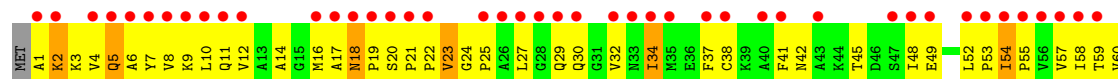
• Molecule 51: 50S RIBOSOMAL PROTEIN L36

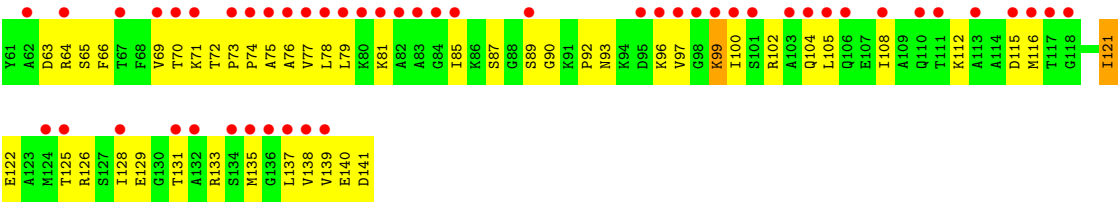


• Molecule 52: 50S RIBOSOMAL PROTEIN L11



• Molecule 52: 50S RIBOSOMAL PROTEIN L11





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 – 3.46 112.47 – 3.53	Depositor EDS
% Data completeness (in resolution range)	91.6 (70.00-3.46) 93.1 (112.47-3.53)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.58Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.279 , 0.331 0.289 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	83.5	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 66.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	284160	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 1.37% 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: KSG, 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	0.23	0/36762	0.71	5/57350 (0.0%)
1	CA	0.23	0/36762	0.72	7/57350 (0.0%)
2	AC	0.23	0/1651	0.44	0/2225
2	CC	0.23	0/1651	0.47	0/2225
3	AD	0.23	0/1665	0.46	0/2227
3	CD	0.23	0/1665	0.45	0/2227
4	AE	0.23	0/1118	0.44	0/1504
4	CE	0.24	0/1118	0.45	0/1504
5	AF	0.24	0/835	0.46	0/1128
5	CF	0.24	0/835	0.47	0/1128
6	AG	0.23	0/1187	0.44	0/1591
6	CG	0.23	0/1211	0.45	0/1624
7	AH	0.23	0/989	0.44	0/1326
7	CH	0.24	0/989	0.45	0/1326
8	AI	0.24	0/1034	0.45	0/1375
8	CI	0.24	0/1034	0.45	0/1375
9	AJ	0.23	0/796	0.49	0/1077
9	CJ	0.22	0/796	0.49	0/1077
10	AK	0.24	0/893	0.46	0/1205
10	CK	0.24	0/893	0.46	0/1205
11	AL	0.22	0/969	0.46	0/1300
11	CL	0.22	0/969	0.47	0/1300
12	AM	0.21	0/892	0.47	0/1193
12	CM	0.21	0/884	0.46	0/1181
13	AN	0.24	0/785	0.45	0/1043
13	CN	0.24	0/785	0.44	0/1043
14	AO	0.23	0/724	0.45	0/966
14	CO	0.23	0/724	0.44	0/966
15	AP	0.26	0/659	0.44	0/884
15	CP	0.25	0/648	0.45	0/870
16	AQ	0.23	0/657	0.46	0/881
16	CQ	0.24	0/666	0.47	0/892

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AR	0.23	0/462	0.45	0/621
17	CR	0.23	0/462	0.45	0/621
18	AS	0.25	0/652	0.47	0/877
18	CS	0.25	0/660	0.46	0/888
19	AT	0.23	0/671	0.41	0/888
19	CT	0.23	0/671	0.42	0/888
20	AB	0.25	0/1735	0.46	0/2338
20	CB	0.25	0/1735	0.46	0/2338
21	AU	0.26	0/430	0.49	0/570
21	CU	0.26	0/430	0.48	0/570
22	BA	0.23	0/2803	0.69	0/4371
22	DA	0.25	0/2803	0.71	0/4371
23	BB	0.26	5/68314 (0.0%)	0.73	22/106569 (0.0%)
23	DB	0.26	5/68314 (0.0%)	0.73	18/106569 (0.0%)
24	BV	0.25	0/766	0.44	0/1025
24	DV	0.25	0/766	0.44	0/1025
25	BC	0.24	0/2092	0.56	0/2813
25	DC	0.23	0/2092	0.56	0/2813
26	BD	0.26	0/1586	0.54	0/2134
26	DD	0.26	0/1586	0.54	0/2134
27	BE	0.25	0/1571	0.54	0/2113
27	DE	0.25	0/1571	0.53	0/2113
28	BF	0.27	0/1444	0.62	0/1937
28	DF	0.26	0/1444	0.59	0/1937
29	BG	0.23	0/1343	0.52	0/1816
29	DG	0.23	0/1343	0.52	0/1816
30	BH	0.25	0/1122	0.51	0/1515
30	DH	0.25	0/1122	0.53	0/1515
31	BJ	0.24	0/1135	0.49	0/1529
31	DJ	0.23	0/1135	0.56	0/1529
32	BK	0.24	0/939	0.63	0/1258
32	DK	0.23	0/939	0.61	0/1258
33	BL	0.27	0/1062	0.72	0/1413
33	DL	0.28	0/1062	0.76	2/1413 (0.1%)
34	BM	0.27	0/1093	0.60	0/1460
34	DM	0.26	0/1093	0.58	0/1460
35	BN	0.24	0/1021	0.52	0/1364
35	DN	0.24	0/1021	0.47	0/1364
36	BO	0.24	0/910	0.50	0/1219
36	DO	0.23	0/910	0.46	0/1219
37	BP	0.26	0/929	0.84	3/1242 (0.2%)
37	DP	0.27	0/929	0.86	3/1242 (0.2%)
38	BQ	0.25	0/960	0.47	0/1278

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	DQ	0.26	0/960	0.50	0/1278
39	BR	0.30	0/829	0.60	0/1107
39	DR	0.26	0/829	0.55	0/1107
40	BS	0.22	0/864	0.50	0/1156
40	DS	0.22	0/864	0.51	0/1156
41	BT	0.24	0/784	0.50	0/1048
41	DT	0.25	0/784	0.54	0/1048
42	BU	0.26	0/787	0.54	0/1051
42	DU	0.25	0/787	0.61	1/1051 (0.1%)
43	BW	0.27	0/642	0.59	0/848
43	DW	0.27	0/642	0.53	0/848
44	BX	0.24	0/510	0.55	0/677
44	DX	0.23	0/510	0.50	0/677
45	BY	0.25	0/453	0.51	0/605
45	DY	0.24	0/453	0.54	0/605
46	BZ	0.28	0/559	0.67	0/745
46	DZ	0.28	0/559	0.61	0/745
47	B0	0.24	0/450	0.52	0/599
47	D0	0.25	0/450	0.58	0/599
48	B1	0.27	0/448	0.51	0/594
48	D1	0.27	0/448	0.50	0/594
49	B2	0.26	0/380	0.48	0/498
49	D2	0.26	0/380	0.47	0/498
50	B3	0.25	0/513	0.54	0/676
50	D3	0.25	0/513	0.51	0/676
51	B4	0.24	0/303	0.54	0/397
51	D4	0.24	0/303	0.53	0/397
52	BI	0.24	0/1046	0.50	0/1410
52	DI	0.26	0/1046	0.53	0/1410
All	All	0.25	10/306470 (0.0%)	0.68	61/458101 (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	14
1	CA	0	16
23	BB	0	41
23	DB	1	41
47	D0	0	1
All	All	1	113

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	BB	1086	A	C5-C6	-17.07	1.25	1.41
23	DB	1086	A	C5-C6	-17.02	1.25	1.41
23	DB	1088	A	C6-N1	-10.55	1.28	1.35
23	BB	1088	A	C6-N1	-10.45	1.28	1.35
23	BB	1060	U	C2-N3	7.88	1.43	1.37
23	DB	1060	U	C2-N3	7.88	1.43	1.37
23	DB	1086	A	N7-C5	-7.07	1.35	1.39
23	BB	1086	A	N7-C5	-7.03	1.35	1.39
23	BB	1086	A	N3-C4	-5.24	1.31	1.34
23	DB	1086	A	N3-C4	-5.11	1.31	1.34

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DB	2791	G	O5'-P-OP2	-26.15	79.33	110.70
23	BB	2791	G	O5'-P-OP1	-26.11	79.37	110.70
23	DB	2791	G	O5'-P-OP1	19.42	134.01	110.70
23	BB	2791	G	O5'-P-OP2	19.11	133.63	110.70
23	DB	2790	U	OP2-P-O3'	15.02	138.25	105.20
23	BB	2790	U	OP1-P-O3'	14.91	138.01	105.20
23	BB	2272	U	C5-C4-O4	-9.03	120.48	125.90
23	DB	2272	U	N3-C4-O4	-8.62	113.36	119.40
23	DB	1088	A	N1-C6-N6	-8.31	113.62	118.60
23	BB	1088	A	N1-C6-N6	-8.21	113.67	118.60
23	BB	1552	A	N9-C1'-C2'	-8.18	103.00	112.00
23	DB	1552	A	N9-C1'-C2'	-8.12	103.07	112.00
23	DB	2076	U	C2'-C3'-O3'	8.12	127.35	109.50
1	CA	765	G	N9-C1'-C2'	-7.70	103.53	112.00
1	AA	765	G	N9-C1'-C2'	-7.67	103.57	112.00
33	DL	118	THR	N-CA-C	7.50	131.26	111.00
23	BB	2733	A	N9-C1'-C2'	-7.46	103.79	112.00
23	DB	2733	A	N9-C1'-C2'	-7.45	103.81	112.00
23	DB	1060	U	C5-C4-O4	-7.35	121.49	125.90
37	BP	71	ARG	N-CA-C	7.26	130.61	111.00
23	BB	1060	U	C5-C4-O4	-7.19	121.59	125.90
37	DP	71	ARG	N-CA-C	7.02	129.96	111.00
23	BB	1086	A	C4-C5-C6	6.85	120.42	117.00
37	BP	79	VAL	N-CA-C	6.84	129.46	111.00
23	DB	1439	A	N9-C1'-C2'	-6.83	104.49	112.00
37	DP	72	VAL	N-CA-C	6.83	129.43	111.00
23	DB	1086	A	C4-C5-C6	6.81	120.40	117.00
23	BB	1439	A	N9-C1'-C2'	-6.76	104.56	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	DB	2790	U	O3'-P-O5'	-6.72	91.22	104.00
33	DL	117	THR	N-CA-C	-6.70	92.92	111.00
1	AA	438	U	N1-C1'-C2'	-6.67	104.66	112.00
1	CA	438	U	N1-C1'-C2'	-6.65	104.69	112.00
1	CA	576	C	O5'-P-OP1	-6.57	99.79	105.70
23	BB	1086	A	C6-C5-N7	-6.52	127.73	132.30
23	DB	1086	A	C6-C5-N7	-6.48	127.76	132.30
37	BP	72	VAL	N-CA-C	6.44	128.39	111.00
23	DB	1088	A	C5-C6-N6	6.41	128.83	123.70
23	BB	1088	A	C5-C6-N6	6.38	128.80	123.70
23	BB	2790	U	O3'-P-O5'	-6.36	91.92	104.00
23	DB	2272	U	C5-C4-O4	-5.80	122.42	125.90
23	BB	1060	U	N1-C2-O2	-5.68	118.82	122.80
23	DB	1060	U	N1-C2-O2	-5.67	118.83	122.80
1	CA	1432	G	N9-C1'-C2'	-5.64	105.79	112.00
23	BB	2267	A	C5-C6-N6	-5.63	119.20	123.70
23	BB	2267	A	O4'-C1'-N9	-5.61	103.72	108.20
23	BB	2267	A	C4-N9-C1'	5.46	136.13	126.30
1	AA	521	G	N9-C1'-C2'	-5.43	106.03	112.00
23	BB	2272	U	N3-C4-O4	-5.41	115.61	119.40
1	CA	521	G	N9-C1'-C2'	-5.41	106.05	112.00
1	CA	328	C	C2'-C3'-O3'	5.39	122.33	113.70
23	DB	1060	U	N3-C2-O2	5.38	125.97	122.20
23	BB	2267	A	C8-N9-C1'	-5.38	118.02	127.70
1	AA	328	C	C2'-C3'-O3'	5.36	122.28	113.70
23	BB	1060	U	N3-C2-O2	5.34	125.94	122.20
23	BB	2272	U	N1-C1'-C2'	-5.33	106.14	112.00
1	CA	1534	A	C2'-C3'-O3'	-5.30	97.84	109.50
23	BB	1086	A	C2-N3-C4	-5.28	107.96	110.60
23	DB	1086	A	C2-N3-C4	-5.16	108.02	110.60
1	AA	66	A	N9-C1'-C2'	-5.09	106.40	112.00
37	DP	70	GLU	N-CA-C	5.07	124.67	111.00
42	DU	49	PRO	N-CA-C	-5.05	98.96	112.10

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
23	DB	2076	U	C3'

All (113) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AA	1047	G	Sidechain
1	AA	1319	A	Sidechain
1	AA	1331	G	Sidechain
1	AA	1441	A	Sidechain
1	AA	187	G	Sidechain
1	AA	281	G	Sidechain
1	AA	324	G	Sidechain
1	AA	437	U	Sidechain
1	AA	438	U	Sidechain
1	AA	454	G	Sidechain
1	AA	496	A	Sidechain
1	AA	521	G	Sidechain
1	AA	575	G	Sidechain
1	AA	703	G	Sidechain
23	BB	1047	G	Sidechain
23	BB	1060	U	Sidechain
23	BB	1086	A	Sidechain
23	BB	1088	A	Sidechain
23	BB	1119	U	Sidechain
23	BB	1240	U	Sidechain
23	BB	1250	G	Sidechain
23	BB	1419	A	Sidechain
23	BB	1439	A	Sidechain
23	BB	1450	G	Sidechain
23	BB	1462	C	Sidechain
23	BB	1546	G	Sidechain
23	BB	1572	A	Sidechain
23	BB	1828	G	Sidechain
23	BB	2108	A	Sidechain
23	BB	221	A	Sidechain
23	BB	2213	U	Sidechain
23	BB	222	A	Sidechain
23	BB	2267	A	Sidechain
23	BB	2272	U	Sidechain
23	BB	2318	G	Sidechain
23	BB	2413	G	Sidechain
23	BB	2471	A	Sidechain
23	BB	2638	G	Sidechain
23	BB	2733	A	Sidechain
23	BB	2770	G	Sidechain
23	BB	2834	G	Sidechain
23	BB	2848	G	Sidechain
23	BB	2857	G	Sidechain

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Mol	Chain	Res	Type	Group
23	BB	2868	A	Sidechain
23	BB	2883	A	Sidechain
23	BB	299	A	Sidechain
23	BB	361	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	729	G	Sidechain
23	BB	757	G	Sidechain
23	BB	942	G	Sidechain
1	CA	1133	G	Sidechain
1	CA	1319	A	Sidechain
1	CA	1362	A	Sidechain
1	CA	1432	G	Sidechain
1	CA	1441	A	Sidechain
1	CA	187	G	Sidechain
1	CA	281	G	Sidechain
1	CA	324	G	Sidechain
1	CA	437	U	Sidechain
1	CA	438	U	Sidechain
1	CA	454	G	Sidechain
1	CA	496	A	Sidechain
1	CA	521	G	Sidechain
1	CA	575	G	Sidechain
1	CA	703	G	Sidechain
1	CA	83	C	Sidechain
47	D0	48	TYR	Sidechain
23	DB	1047	G	Sidechain
23	DB	1060	U	Sidechain
23	DB	1086	A	Sidechain
23	DB	1088	A	Sidechain
23	DB	1247	A	Sidechain
23	DB	136	G	Sidechain
23	DB	1419	A	Sidechain
23	DB	1439	A	Sidechain
23	DB	1450	G	Sidechain
23	DB	1462	C	Sidechain
23	DB	1546	G	Sidechain
23	DB	1572	A	Sidechain
23	DB	1828	G	Sidechain

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Mol	Chain	Res	Type	Group
23	DB	2062	A	Sidechain
23	DB	2108	A	Sidechain
23	DB	214	G	Sidechain
23	DB	221	A	Sidechain
23	DB	222	A	Sidechain
23	DB	2267	A	Sidechain
23	DB	2272	U	Sidechain
23	DB	232	G	Sidechain
23	DB	2413	G	Sidechain
23	DB	2471	A	Sidechain
23	DB	2503	A	Sidechain
23	DB	2638	G	Sidechain
23	DB	2733	A	Sidechain
23	DB	2770	G	Sidechain
23	DB	28	A	Sidechain
23	DB	2834	G	Sidechain
23	DB	2848	G	Sidechain
23	DB	2857	G	Sidechain
23	DB	2868	A	Sidechain
23	DB	2883	A	Sidechain
23	DB	299	A	Sidechain
23	DB	500	G	Sidechain
23	DB	630	G	Sidechain
23	DB	633	A	Sidechain
23	DB	727	A	Sidechain
23	DB	729	G	Sidechain
23	DB	757	G	Sidechain
23	DB	858	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	1330	0
1	CA	32831	0	16521	1382	0
2	AC	1624	0	1699	184	0
2	CC	1624	0	1699	162	0
3	AD	1643	0	1710	173	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	CD	1643	0	1710	175	0
4	AE	1105	0	1148	108	0
4	CE	1105	0	1148	143	0
5	AF	817	0	808	73	0
5	CF	817	0	808	102	0
6	AG	1174	0	1230	104	0
6	CG	1196	0	1246	101	0
7	AH	979	0	1034	85	0
7	CH	979	0	1034	80	0
8	AI	1022	0	1070	156	0
8	CI	1022	0	1070	125	0
9	AJ	786	0	828	95	0
9	CJ	786	0	828	111	0
10	AK	877	0	887	104	0
10	CK	877	0	887	106	0
11	AL	955	0	1019	105	0
11	CL	955	0	1019	91	0
12	AM	883	0	944	88	0
12	CM	876	0	937	113	0
13	AN	774	0	827	104	0
13	CN	774	0	827	121	0
14	AO	716	0	742	53	0
14	CO	716	0	742	55	0
15	AP	649	0	666	78	0
15	CP	638	0	656	79	0
16	AQ	648	0	691	101	0
16	CQ	657	0	702	90	0
17	AR	455	0	478	42	0
17	CR	455	0	478	45	0
18	AS	637	0	665	87	0
18	CS	644	0	675	96	0
19	AT	665	0	714	59	0
19	CT	665	0	714	58	0
20	AB	1704	0	1732	195	0
20	CB	1704	0	1732	167	0
21	AU	425	0	449	68	0
21	CU	425	0	449	65	0
22	BA	2507	0	1270	84	0
22	DA	2507	0	1270	117	0
23	BB	60995	0	30678	2753	0
23	DB	60995	0	30677	2695	0
24	BV	753	0	780	100	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
24	DV	753	0	780	69	0
25	BC	2053	0	2122	451	0
25	DC	2053	0	2122	389	0
26	BD	1565	0	1616	379	0
26	DD	1565	0	1616	309	0
27	BE	1552	0	1619	279	0
27	DE	1552	0	1619	256	0
28	BF	1420	0	1460	170	0
28	DF	1420	0	1460	157	0
29	BG	1323	0	1374	166	0
29	DG	1323	0	1374	158	0
30	BH	1111	0	1148	173	0
30	DH	1111	0	1148	166	0
31	BJ	1112	0	1147	220	0
31	DJ	1112	0	1147	207	0
32	BK	930	0	1000	123	0
32	DK	930	0	1000	124	0
33	BL	1053	0	1129	277	0
33	DL	1053	0	1129	250	0
34	BM	1074	0	1157	238	0
34	DM	1074	0	1157	169	0
35	BN	1008	0	1045	186	0
35	DN	1008	0	1045	158	0
36	BO	900	0	935	128	0
36	DO	900	0	935	137	0
37	BP	917	0	965	198	0
37	DP	917	0	965	196	0
38	BQ	947	0	1022	199	0
38	DQ	947	0	1022	175	0
39	BR	816	0	839	170	0
39	DR	816	0	839	166	0
40	BS	857	0	922	124	0
40	DS	857	0	922	116	0
41	BT	777	0	840	148	0
41	DT	777	0	840	133	0
42	BU	779	0	834	163	0
42	DU	779	0	834	121	0
43	BW	634	0	656	161	0
43	DW	634	0	656	169	0
44	BX	509	0	543	77	0
44	DX	509	0	543	83	0
45	BY	449	0	491	60	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
45	DY	449	0	491	81	0
46	BZ	549	0	552	111	0
46	DZ	549	0	552	102	0
47	B0	444	0	461	74	0
47	D0	444	0	461	91	0
48	B1	441	0	485	80	0
48	D1	441	0	485	77	0
49	B2	377	0	418	59	0
49	D2	377	0	418	67	0
50	B3	504	0	574	115	0
50	D3	504	0	574	107	0
51	B4	302	0	343	63	0
51	D4	302	0	343	75	0
52	BI	1032	0	1088	121	0
52	DI	1032	0	1088	183	0
53	AA	26	0	23	3	0
53	CA	26	0	23	1	0
54	AA	60	0	0	0	0
54	BB	110	0	0	0	0
54	CA	62	0	0	0	0
54	DB	109	0	0	0	0
54	DE	1	0	0	0	0
54	DN	1	0	0	0	0
55	AA	289	0	0	1	0
55	AE	3	0	0	0	0
55	AK	2	0	0	0	0
55	AN	3	0	0	0	0
55	AP	2	0	0	0	0
55	AT	1	0	0	0	0
55	BB	497	0	0	11	0
55	BC	1	0	0	0	0
55	BE	5	0	0	0	0
55	BH	1	0	0	0	0
55	BL	2	0	0	0	0
55	BN	1	0	0	0	0
55	CA	293	0	0	1	0
55	CE	3	0	0	0	0
55	CK	1	0	0	0	0
55	CL	4	0	0	0	0
55	CN	3	0	0	0	0
55	CP	1	0	0	0	0
55	CT	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
55	D2	2	0	0	0	0
55	DB	501	0	0	10	0
55	DC	1	0	0	0	0
55	DD	1	0	0	0	0
55	DE	3	0	0	0	0
55	DL	1	0	0	0	0
55	DN	2	0	0	0	0
55	DT	1	0	0	0	0
All	All	284160	0	190815	19652	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 42.

All (19652) 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:O5'	52:DI:4:VAL:N	1.71	1.23
32:DK:78:ARG:HG2	37:DP:72:VAL:HG21	1.24	1.15
23:DB:1098:A:H3'	52:DI:3:LYS:CA	1.76	1.15
23:DB:587:C:H3'	33:DL:29:LYS:HD2	1.20	1.14
48:D1:29:LYS:HB2	48:D1:30:PRO:HD3	1.30	1.14
1:AA:699:C:H2'	1:AA:700:G:H5''	1.28	1.14
23:BB:662:G:H4'	33:BL:24:GLY:H	1.11	1.13
42:DU:60:LYS:HE3	42:DU:61:GLU:H	1.08	1.13
23:DB:458:G:H5''	49:D2:39:ARG:HB2	1.30	1.12
44:DX:43:LEU:HB2	44:DX:45:GLN:HE22	1.13	1.12
23:BB:85:G:H5'	42:BU:28:LEU:HA	1.30	1.11
1:CA:243:A:H4'	1:CA:244:U:H5'	1.29	1.11
20:AB:163:ILE:HG23	20:AB:164:ASP:H	1.14	1.11
33:DL:7:SER:HB2	33:DL:8:PRO:HD3	1.24	1.11
36:BO:57:ALA:HB2	36:BO:62:LEU:HD22	1.30	1.10
32:BK:108:ARG:HH21	37:BP:36:LYS:HG3	1.06	1.10
30:BH:82:SER:H	30:BH:146:VAL:HG13	1.16	1.10
23:BB:2750:A:H4'	29:BG:1:SER:HB2	1.34	1.09
34:BM:3:GLN:H	34:BM:4:PRO:HD2	1.04	1.09
34:DM:57:VAL:HG21	34:DM:112:LEU:HD21	1.33	1.09
23:BB:38:A:O4'	27:BE:46:GLN:HG3	1.53	1.08
20:CB:46:VAL:HG13	20:CB:47:PRO:HD3	1.35	1.08
25:DC:33:LEU:HD22	25:DC:34:GLU:HG3	1.35	1.08
11:CL:43:LYS:HB3	11:CL:44:PRO:HD2	1.28	1.08
1:CA:974:A:H4'	1:CA:975:A:H5'	1.28	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:699:C:H2'	1:CA:700:G:H5''	1.28	1.08
23:DB:2527:C:H5''	51:D4:34:LYS:HG3	1.35	1.08
1:AA:243:A:H4'	1:AA:244:U:H5'	1.29	1.07
26:BD:24:VAL:HG21	26:BD:193:VAL:HG11	1.36	1.06
43:DW:38:ARG:HH21	43:DW:40:ARG:HD3	1.16	1.06
25:BC:42:ARG:HG2	25:BC:43:ASN:H	1.14	1.06
23:BB:100:U:H4'	42:BU:90:LYS:HE2	1.36	1.05
23:DB:1098:A:H3'	52:DI:3:LYS:HA	1.37	1.05
25:BC:107:LYS:HB2	25:BC:194:VAL:HG11	1.39	1.05
23:BB:161:A:H3'	23:BB:162:U:H5''	1.35	1.05
23:BB:2204:G:H5'	25:BC:149:LYS:HE3	1.36	1.04
23:DB:45:G:H5''	23:DB:46:G:H5'	1.36	1.04
23:BB:2333:A:H4'	23:BB:2334:U:H5''	1.40	1.04
30:DH:3:VAL:HG22	30:DH:21:VAL:HG11	1.39	1.04
33:BL:14:LYS:HA	33:BL:14:LYS:HZ3	1.20	1.04
25:DC:107:LYS:HB2	25:DC:194:VAL:HG21	1.38	1.04
23:DB:1099:G:H8	52:DI:3:LYS:N	1.56	1.04
35:DN:45:ARG:HH22	35:DN:113:ILE:HG23	1.21	1.04
31:DJ:102:GLU:HG3	31:DJ:124:VAL:HG12	1.40	1.03
31:BJ:135:GLN:HG2	31:BJ:137:PRO:HD2	1.40	1.03
43:BW:47:GLY:HA3	43:BW:55:ASP:HB3	1.39	1.03
37:BP:52:ARG:HB3	37:BP:60:VAL:HG11	1.37	1.03
38:DQ:48:ASP:HA	38:DQ:51:GLN:HE21	1.20	1.03
18:AS:18:VAL:HG21	18:AS:43:MET:HG2	1.35	1.03
36:DO:56:LYS:HE2	36:DO:81:ARG:HE	1.16	1.03
27:BE:128:ALA:HB3	27:BE:129:PRO:HD3	1.36	1.03
20:AB:46:VAL:HG13	20:AB:47:PRO:HD3	1.38	1.03
30:BH:37:VAL:H	30:BH:38:PRO:HD2	1.23	1.03
23:BB:923:G:H1'	43:BW:23:LYS:HD3	1.38	1.03
42:DU:28:LEU:HD13	42:DU:30:SER:H	1.24	1.03
23:DB:323:C:H1'	27:DE:164:LEU:HB3	1.39	1.02
33:DL:51:GLU:HG2	33:DL:52:GLY:H	1.24	1.02
51:D4:26:ILE:HG13	51:D4:35:GLN:H	1.22	1.02
30:DH:6:LEU:HD12	30:DH:36:ALA:H	1.24	1.02
37:DP:93:LYS:HG2	37:DP:96:LEU:HA	1.41	1.02
33:BL:61:LEU:HB3	33:BL:62:PRO:HD2	1.42	1.02
15:AP:42:ILE:HB	15:AP:46:LYS:HD2	1.42	1.02
32:BK:71:ARG:HB3	32:BK:72:PRO:HD2	1.04	1.02
45:BY:16:LEU:HB2	45:BY:17:PRO:HD3	1.37	1.02
33:BL:13:LYS:HG3	33:BL:14:LYS:H	1.25	1.02
46:BZ:8:LYS:HD2	46:BZ:9:TYR:H	1.22	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:BK:71:ARG:HB3	32:BK:72:PRO:CD	1.90	1.01
32:BK:8:LEU:HD23	32:BK:82:ASN:HB3	1.41	1.01
23:DB:1099:G:P	52:DI:4:VAL:H	1.83	1.01
37:DP:47:ILE:HG22	37:DP:48:ALA:H	1.22	1.01
23:DB:2091:C:H3'	23:DB:2092:U:H5''	1.43	1.01
8:CI:27:ILE:HG23	8:CI:34:LEU:HB2	1.42	1.00
48:D1:46:VAL:HG13	48:D1:47:ILE:HG13	1.41	1.00
23:DB:2333:A:H4'	23:DB:2334:U:H5''	1.40	1.00
34:BM:38:ARG:HB2	34:BM:38:ARG:HH11	1.26	1.00
23:BB:2052:A:H4'	26:BD:148:GLN:H	1.24	1.00
23:BB:1454:C:H5'	35:BN:63:ARG:HH21	1.24	1.00
23:DB:1025:G:H1'	23:DB:1135:C:H5'	1.43	1.00
24:DV:63:ILE:H	24:DV:70:ILE:HD11	1.23	1.00
23:BB:442:G:H1'	27:BE:44:ARG:HD2	1.38	1.00
23:BB:692:C:H5'	25:BC:42:ARG:HB2	1.40	1.00
16:AQ:76:ARG:HE	16:AQ:78:VAL:HG22	1.27	1.00
11:AL:8:ARG:HG3	11:AL:9:LYS:H	1.24	1.00
5:CF:3:HIS:HB3	5:CF:92:THR:HA	1.41	1.00
23:DB:136:G:H2'	23:DB:137:U:C6	1.97	0.99
20:AB:67:LEU:HD11	20:AB:157:PRO:HB3	1.44	0.99
25:DC:107:LYS:HE3	25:DC:108:GLY:H	1.23	0.99
26:DD:197:THR:HG23	26:DD:198:GLY:H	1.28	0.99
46:DZ:55:GLY:HA2	46:DZ:59:ARG:HB2	1.43	0.99
3:AD:7:LYS:HE2	3:AD:20:LEU:HD13	1.43	0.99
25:DC:230:PRO:HG2	25:DC:245:THR:H	1.27	0.99
25:DC:171:VAL:HB	25:DC:182:LYS:HB3	1.45	0.99
1:CA:973:G:H3'	1:CA:974:A:H5''	1.45	0.99
13:AN:60:ARG:HD3	13:AN:60:ARG:H	1.28	0.98
25:BC:139:THR:HG22	25:BC:193:GLU:HB3	1.44	0.98
43:DW:13:ARG:HE	43:DW:13:ARG:H	1.08	0.98
23:DB:2377:A:H61	36:DO:13:ARG:HH21	1.07	0.98
9:AJ:35:GLN:HB3	9:AJ:77:VAL:HG23	1.40	0.98
27:DE:169:VAL:HG13	27:DE:170:ARG:H	1.29	0.98
41:DT:66:LYS:H	41:DT:76:ARG:HH21	1.11	0.98
23:DB:858:G:N3	23:DB:2268:A:H2'	1.78	0.98
39:BR:78:ARG:HB3	39:BR:87:GLN:HA	1.45	0.97
25:DC:48:ILE:HG22	25:DC:49:THR:H	1.26	0.97
23:BB:2266:A:H4'	23:BB:2267:A:N7	1.80	0.97
50:D3:49:VAL:HG22	50:D3:50:SER:H	1.28	0.97
51:D4:11:CYS:HB2	51:D4:14:CYS:HB2	1.46	0.97
33:BL:29:LYS:HD2	33:BL:31:GLY:H	1.27	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BS:6:LYS:HA	40:BS:104:THR:HA	1.46	0.97
23:BB:2478:A:H5'	51:B4:32:LYS:HD2	1.47	0.97
45:BY:10:ARG:HE	45:BY:31:ILE:HG13	1.27	0.97
23:BB:396:G:H4'	46:BZ:28:VAL:HG21	1.46	0.97
25:BC:53:ILE:HD13	25:BC:218:THR:HG23	1.45	0.97
46:BZ:33:ASN:HB2	46:BZ:44:PHE:HB2	1.44	0.97
46:BZ:3:LYS:HB2	46:BZ:51:VAL:HG21	1.45	0.97
23:DB:1099:G:C8	52:DI:3:LYS:N	2.32	0.97
39:DR:63:VAL:HG22	39:DR:64:VAL:H	1.28	0.97
23:BB:482:A:H5''	42:BU:43:LYS:HE3	1.47	0.97
32:BK:26:GLY:HA3	32:BK:30:ARG:HE	1.29	0.97
40:DS:46:LEU:HA	40:DS:49:LYS:HB2	1.45	0.97
41:DT:48:GLN:HA	41:DT:53:VAL:HG22	1.46	0.97
23:BB:1450:G:H21	23:BB:1452:G:H1	1.12	0.97
23:BB:528:A:N1	23:BB:2042:A:H2'	1.80	0.97
26:BD:170:VAL:HG13	26:BD:171:THR:H	1.29	0.96
32:BK:71:ARG:CB	32:BK:72:PRO:HD2	1.95	0.96
22:BA:75:G:H1'	24:BV:29:ILE:HD13	1.44	0.96
8:AI:29:ILE:HG22	8:AI:64:ILE:HB	1.47	0.96
23:BB:858:G:H21	23:BB:2268:A:H3'	1.31	0.96
31:DJ:35:ARG:HH12	31:DJ:40:HIS:H	0.98	0.96
45:DY:18:LYS:HD2	45:DY:18:LYS:H	1.28	0.96
38:DQ:53:LYS:HE2	38:DQ:53:LYS:H	1.30	0.96
28:BF:36:ASN:HB2	28:BF:87:LYS:HA	1.47	0.96
31:BJ:81:ILE:HG13	31:BJ:82:GLY:H	1.26	0.96
35:BN:113:ILE:HG22	35:BN:114:GLU:H	1.30	0.96
1:CA:1320:C:H5''	18:CS:2:ARG:HD3	1.47	0.96
31:DJ:124:VAL:HG23	31:DJ:125:TYR:H	1.27	0.96
21:CU:17:ARG:H	21:CU:17:ARG:HD2	1.29	0.96
51:D4:3:VAL:HG12	51:D4:4:ARG:H	1.30	0.96
25:BC:28:PRO:HD2	25:BC:79:ARG:HH21	1.27	0.96
1:CA:1003:G:H21	1:CA:1005:A:H5'	1.27	0.96
39:BR:74:ILE:HA	39:BR:90:ARG:HE	1.27	0.96
23:BB:858:G:N3	23:BB:2268:A:H2'	1.81	0.96
1:CA:1151:A:HO2'	1:CA:1152:A:H8	1.12	0.96
10:CK:78:ILE:HD13	10:CK:78:ILE:H	1.28	0.96
34:DM:5:LYS:HE3	34:DM:6:ARG:N	1.80	0.96
37:DP:25:VAL:HG11	37:DP:87:ARG:HA	1.47	0.96
42:DU:66:VAL:HG22	42:DU:67:SER:H	1.29	0.96
13:AN:63:CYS:HB3	13:AN:67:GLY:H	1.30	0.95
38:BQ:92:LYS:HG2	38:BQ:93:ILE:H	1.29	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:29:ILE:HD13	5:CF:64:VAL:HG11	1.48	0.95
46:DZ:54:GLY:H	46:DZ:57:VAL:HG23	1.31	0.95
31:DJ:40:HIS:HB2	38:DQ:69:ARG:HH22	1.29	0.95
20:CB:156:LEU:HD11	20:CB:178:LEU:HD13	1.47	0.95
25:BC:28:PRO:HD2	25:BC:79:ARG:NH2	1.82	0.95
23:DB:1368:G:H5''	49:D2:25:LYS:HG2	1.46	0.95
23:BB:1392:A:H62	41:BT:18:GLU:HG2	1.30	0.95
41:BT:88:LYS:HG2	41:BT:89:GLU:H	1.30	0.95
19:AT:68:LYS:HG3	19:AT:69:ASN:H	1.29	0.95
23:BB:2576:G:P	26:BD:149:ASN:HB2	2.07	0.95
31:DJ:84:ILE:HD12	31:DJ:85:LYS:H	1.31	0.95
27:DE:127:GLU:HA	27:DE:157:LEU:HD13	1.49	0.95
5:AF:53:LYS:HD3	5:AF:54:LEU:H	1.30	0.94
18:AS:49:ALA:HB1	18:AS:56:HIS:HB3	1.49	0.94
12:AM:48:SER:H	12:AM:51:GLN:HB2	1.33	0.94
1:CA:437:U:H2'	1:CA:438:U:O4'	1.67	0.94
40:BS:71:VAL:HG22	40:BS:107:VAL:HG12	1.45	0.94
44:BX:24:GLU:HA	44:BX:28:LEU:HB2	1.47	0.94
23:DB:2377:A:H61	36:DO:13:ARG:NH2	1.64	0.94
46:DZ:30:HIS:HB2	46:DZ:48:GLN:HG2	1.48	0.94
1:AA:764:C:H3'	1:AA:765:G:H21	1.33	0.94
27:BE:116:ASP:HB2	27:BE:185:LYS:HA	1.50	0.94
43:BW:35:ILE:HG23	43:BW:36:ILE:H	1.32	0.94
8:CI:118:ARG:HH21	8:CI:122:ARG:HE	1.13	0.94
39:DR:65:ALA:HB3	39:DR:99:THR:HG23	1.48	0.94
27:BE:47:LYS:HD3	27:BE:47:LYS:H	1.30	0.94
38:BQ:93:ILE:HG23	38:BQ:94:LEU:HD13	1.50	0.94
46:BZ:36:VAL:HA	46:BZ:42:PRO:HB3	1.45	0.94
3:CD:167:PRO:HG2	3:CD:170:LEU:HD11	1.48	0.94
23:DB:918:A:H2'	23:DB:919:U:H5'	1.46	0.94
21:AU:3:ILE:HA	21:AU:19:LYS:HG2	1.50	0.94
51:B4:12:ARG:HD2	51:B4:12:ARG:H	1.32	0.94
10:CK:19:VAL:HG22	10:CK:34:THR:HG22	1.49	0.94
34:DM:57:VAL:HG12	34:DM:58:LYS:H	1.30	0.94
1:AA:437:U:H2'	1:AA:438:U:O4'	1.67	0.94
28:BF:133:GLU:HG2	28:BF:147:ARG:HG2	1.49	0.94
37:DP:47:ILE:HG23	37:DP:63:ILE:HA	1.46	0.94
24:BV:4:ILE:HB	24:BV:63:ILE:HA	1.49	0.93
23:DB:972:A:H3'	23:DB:973:A:H5''	1.48	0.93
23:DB:1199:U:H2'	23:DB:1200:C:H6	1.31	0.93
31:DJ:41:LYS:HZ1	38:DQ:63:ARG:HD2	1.33	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1313:U:H5''	18:AS:5:LYS:HG2	1.49	0.93
13:CN:50:LEU:H	13:CN:51:PRO:HD2	1.33	0.93
23:DB:1450:G:H21	23:DB:1452:G:H1	1.12	0.93
26:DD:77:ARG:HH21	26:DD:79:LEU:HB2	1.30	0.93
30:DH:125:THR:HA	30:DH:146:VAL:HB	1.51	0.93
41:BT:54:GLU:H	41:BT:91:GLN:HE22	1.15	0.93
23:DB:142:A:H2'	23:DB:143:C:C6	2.03	0.93
28:DF:28:PRO:HB2	28:DF:168:LEU:HD12	1.48	0.93
44:BX:28:LEU:HB3	44:BX:34:SER:HA	1.49	0.93
1:CA:264:C:H4'	16:CQ:64:ARG:HD2	1.49	0.93
37:DP:26:GLU:HA	37:DP:47:ILE:H	1.33	0.93
12:AM:78:ARG:HH21	12:AM:79:LEU:HG	1.33	0.93
23:DB:2574:G:H21	26:DD:147:GLY:HA3	1.31	0.93
23:DB:2304:G:H4'	28:DF:129:MET:HA	1.48	0.93
38:BQ:33:VAL:HG13	38:BQ:34:ALA:H	1.31	0.93
28:DF:32:LYS:HB3	28:DF:91:ARG:HB3	1.48	0.93
10:AK:22:ILE:HG21	10:AK:95:THR:HG21	1.51	0.93
14:AO:6:ALA:HA	14:AO:9:LYS:HE3	1.50	0.93
33:DL:85:VAL:HG13	33:DL:86:GLU:H	1.33	0.93
37:DP:100:ARG:HH11	37:DP:100:ARG:HB3	1.32	0.93
23:BB:37:C:H2'	27:BE:46:GLN:CG	1.99	0.93
26:BD:16:THR:HG23	26:BD:17:GLU:H	1.33	0.93
33:BL:14:LYS:NZ	33:BL:14:LYS:HA	1.82	0.93
33:BL:94:THR:HG23	33:BL:95:LEU:HD12	1.50	0.93
49:B2:42:LEU:HD11	49:B2:44:VAL:HG13	1.48	0.92
23:BB:919:U:H2'	23:BB:920:A:C8	2.05	0.92
41:BT:8:LEU:HB3	44:BX:19:LEU:HD11	1.48	0.92
43:DW:35:ILE:HD12	43:DW:35:ILE:H	1.34	0.92
23:BB:480:A:H4'	42:BU:42:LYS:HG2	1.51	0.92
32:BK:66:LYS:HG2	32:BK:80:ASP:HA	1.49	0.92
46:BZ:25:ARG:HB2	46:BZ:25:ARG:HH11	1.33	0.92
23:DB:2046:G:H5'	47:D0:15:ARG:HD2	1.51	0.92
23:DB:1099:G:P	52:DI:3:LYS:HA	2.09	0.92
23:BB:535:G:H21	38:BQ:52:ARG:HH22	1.13	0.92
27:BE:48:THR:HG22	27:BE:49:ARG:H	1.34	0.92
33:BL:60:ARG:HA	33:BL:60:ARG:HH11	1.31	0.92
48:D1:47:ILE:HG22	48:D1:48:TYR:H	1.34	0.92
22:DA:48:U:H2'	22:DA:49:C:C6	2.03	0.92
33:BL:90:VAL:HG11	33:BL:121:THR:H	1.34	0.92
21:CU:24:LYS:HD2	21:CU:25:ALA:H	1.34	0.92
23:DB:1024:G:H3'	23:DB:1025:G:H5''	1.50	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DF:13:LYS:HA	28:DF:16:MET:HB2	1.51	0.92
32:BK:70:ARG:HB3	32:BK:76:VAL:HG22	1.50	0.92
39:BR:47:VAL:HG13	39:BR:49:ILE:H	1.34	0.92
12:CM:52:ILE:HG13	12:CM:56:ARG:HH22	1.33	0.92
23:DB:1099:G:H8	52:DI:3:LYS:CA	1.83	0.92
41:DT:76:ARG:HH12	41:DT:78:SER:HB2	1.35	0.92
27:BE:143:LEU:HG	27:BE:185:LYS:HD2	1.51	0.92
30:BH:134:VAL:HG13	30:BH:135:HIS:H	1.35	0.92
26:BD:118:PHE:HA	26:BD:164:GLN:HA	1.52	0.92
27:DE:142:ALA:H	27:DE:185:LYS:HZ1	1.12	0.92
27:DE:148:ILE:HA	27:DE:185:LYS:HB3	1.50	0.92
37:DP:27:VAL:HG13	37:DP:47:ILE:HD12	1.49	0.92
1:AA:1086:U:H3	1:AA:1099:G:H22	1.05	0.92
23:BB:543:G:H2'	23:BB:544:C:H5''	1.51	0.92
28:BF:168:LEU:HD12	28:BF:169:LEU:H	1.35	0.92
3:CD:58:GLN:HA	3:CD:58:GLN:HE21	1.35	0.92
9:CJ:51:VAL:HG23	13:CN:80:ARG:HB2	1.51	0.92
26:BD:126:ASN:O	26:BD:127:PHE:HB2	1.70	0.92
33:BL:77:ILE:HB	33:BL:110:VAL:HG22	1.51	0.92
1:CA:1367:C:H5''	8:CI:115:VAL:HG23	1.49	0.92
31:BJ:64:VAL:HG11	31:BJ:68:LYS:HD2	1.50	0.91
1:AA:376:G:H5''	15:AP:5:ARG:HB2	1.52	0.91
10:AK:124:LYS:HA	21:AU:34:ARG:HB3	1.49	0.91
23:BB:630:G:H1	33:BL:69:ARG:HH22	1.04	0.91
34:BM:96:ILE:HG23	34:BM:97:GLN:H	1.35	0.91
21:CU:4:LYS:HB3	21:CU:6:ARG:HH12	1.32	0.91
40:DS:23:LEU:HB2	47:D0:21:LEU:HD13	1.52	0.91
23:DB:2054:A:H2'	47:D0:4:GLN:HE22	1.35	0.91
30:DH:26:ALA:HB2	30:DH:30:LEU:HG	1.49	0.91
23:DB:856:G:H4'	43:DW:23:LYS:HD2	1.52	0.91
20:AB:9:LEU:HD22	20:AB:11:ALA:H	1.33	0.91
23:BB:1639:C:H2'	23:BB:1640:A:H5''	1.52	0.91
33:BL:55:MET:HG2	33:BL:56:PRO:HD3	1.52	0.91
23:DB:27:G:H22	23:DB:512:G:H2'	1.35	0.91
25:DC:127:ASN:HD22	25:DC:128:THR:H	1.15	0.91
31:BJ:45:THR:HB	38:BQ:63:ARG:HH22	1.35	0.91
36:DO:40:ILE:HD13	36:DO:40:ILE:H	1.35	0.91
41:DT:55:VAL:HG22	41:DT:56:GLU:H	1.33	0.91
10:AK:86:LYS:HG3	10:AK:113:THR:HA	1.53	0.91
23:BB:1453:A:N6	35:BN:74:GLU:HG2	1.85	0.91
38:BQ:91:ARG:H	39:BR:10:LYS:HZ2	0.93	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:360:U:H2'	23:DB:361:G:O4'	1.70	0.91
23:BB:670:A:H3'	33:BL:47:ARG:HH11	1.32	0.91
27:DE:4:VAL:HG13	27:DE:5:LEU:H	1.36	0.91
10:CK:20:ALA:HB3	10:CK:83:VAL:HA	1.53	0.91
30:BH:5:LEU:H	30:BH:5:LEU:HD22	1.35	0.91
1:CA:764:C:H3'	1:CA:765:G:H21	1.33	0.91
2:CC:59:PRO:HG2	2:CC:62:SER:HB2	1.49	0.91
50:D3:12:ARG:HB3	50:D3:23:HIS:HB2	1.52	0.91
31:DJ:81:ILE:HG23	31:DJ:82:GLY:H	1.34	0.91
40:DS:29:VAL:HG22	40:DS:71:VAL:HG23	1.50	0.91
50:B3:12:ARG:HB3	50:B3:23:HIS:HA	1.53	0.91
36:DO:66:GLY:H	36:DO:70:ALA:HB2	1.34	0.91
6:AG:125:ASP:HB3	6:AG:130:LYS:HB3	1.53	0.90
37:DP:47:ILE:HG12	37:DP:63:ILE:HG23	1.52	0.90
26:DD:29:VAL:HG22	26:DD:30:GLU:H	1.36	0.90
26:BD:115:GLY:HA2	26:BD:165:MET:HG3	1.53	0.90
51:D4:18:LYS:HE2	51:D4:19:ARG:N	1.86	0.90
21:AU:13:VAL:HG13	21:AU:14:ALA:H	1.35	0.90
49:B2:44:VAL:HG23	49:B2:45:SER:H	1.37	0.90
27:BE:60:TRP:HE1	27:BE:73:ILE:HD11	1.34	0.90
23:DB:1060:U:N3	23:DB:1088:A:N7	2.20	0.90
32:DK:78:ARG:HH22	37:DP:62:LYS:HZ2	1.16	0.90
25:BC:30:ALA:H	25:BC:31:PRO:HD3	1.36	0.90
23:BB:691:C:O2'	25:BC:42:ARG:HD2	1.70	0.90
23:BB:811:U:H2'	33:BL:31:GLY:HA3	1.53	0.90
38:BQ:5:ARG:HB2	38:BQ:5:ARG:HH11	1.34	0.90
25:DC:224:MET:HA	25:DC:233:GLY:H	1.36	0.90
2:AC:156:LEU:H	2:AC:156:LEU:HD12	1.33	0.90
25:BC:143:VAL:HG12	25:BC:144:GLU:H	1.35	0.90
1:CA:451:A:H5'	15:CP:70:ARG:HH22	1.35	0.90
1:CA:719:C:H1'	17:CR:37:LYS:HB2	1.52	0.90
23:DB:2645:G:H3'	23:DB:2646:C:H5'	1.54	0.90
28:DF:140:ILE:HD12	28:DF:140:ILE:H	1.37	0.90
37:BP:50:ARG:HH11	37:BP:50:ARG:HB2	1.33	0.90
18:CS:29:PRO:HA	18:CS:47:THR:HB	1.54	0.90
50:B3:24:LYS:HA	50:B3:45:PRO:HB3	1.54	0.90
20:CB:15:PHE:HA	20:CB:42:LEU:HD21	1.51	0.90
50:D3:32:LEU:HD13	50:D3:33:THR:H	1.35	0.90
27:DE:149:ILE:HD13	27:DE:186:VAL:HG13	1.54	0.90
32:BK:103:VAL:HB	32:BK:107:LEU:HD23	1.54	0.90
41:DT:31:VAL:HG13	41:DT:32:LEU:H	1.37	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1211:U:H4'	1:AA:1213:A:H1'	1.54	0.89
20:AB:202:ASN:HD22	20:AB:204:ASP:H	1.20	0.89
23:BB:1283:G:H22	23:BB:1286:A:H5'	1.36	0.89
34:BM:3:GLN:N	34:BM:4:PRO:HD2	1.82	0.89
40:BS:27:LYS:H	40:BS:27:LYS:HD2	1.33	0.89
5:CF:38:ARG:HH11	5:CF:98:GLU:H	1.18	0.89
31:DJ:98:GLU:HB3	31:DJ:124:VAL:HB	1.51	0.89
7:AH:103:VAL:HG12	7:AH:124:ILE:HA	1.53	0.89
23:BB:2575:C:H4'	26:BD:148:GLN:HA	1.54	0.89
31:BJ:78:THR:H	31:BJ:84:ILE:HG21	1.35	0.89
34:BM:70:ASP:N	34:BM:71:LYS:HZ2	1.69	0.89
45:BY:9:THR:HA	45:BY:55:LYS:HE2	1.53	0.89
31:DJ:58:ASN:O	31:DJ:126:ALA:HA	1.71	0.89
44:DX:18:LEU:HD22	44:DX:18:LEU:H	1.36	0.89
23:BB:2645:G:H3'	23:BB:2646:C:H5'	1.54	0.89
23:BB:855:G:N2	43:BW:23:LYS:HD2	1.86	0.89
23:BB:45:G:H5'	23:BB:46:G:H5'	1.53	0.89
26:BD:2:ILE:HG21	26:BD:204:LYS:HA	1.52	0.89
52:BI:27:LEU:HD23	52:BI:27:LEU:H	1.37	0.89
41:BT:58:VAL:HG13	41:BT:85:VAL:HG12	1.51	0.89
23:DB:161:A:H3'	23:DB:162:U:H5''	1.53	0.89
24:DV:25:LYS:HE2	24:DV:41:GLU:HB2	1.54	0.89
5:CF:6:ILE:HG23	5:CF:62:MET:HB3	1.52	0.89
34:DM:18:ARG:HD3	34:DM:18:ARG:H	1.36	0.89
46:BZ:5:ILE:HD12	46:BZ:5:ILE:H	1.37	0.89
25:DC:50:THR:HG22	25:DC:51:ARG:HG3	1.52	0.89
25:BC:229:HIS:NE2	25:BC:244:VAL:HG22	1.88	0.89
33:BL:77:ILE:HD13	33:BL:92:LEU:HD13	1.52	0.89
23:DB:1283:G:H22	23:DB:1286:A:H5'	1.36	0.89
30:DH:86:ASP:HB2	30:DH:89:LYS:HD3	1.55	0.89
23:BB:1060:U:N3	23:BB:1088:A:N7	2.20	0.89
33:BL:30:THR:HB	33:BL:36:LYS:HB2	1.55	0.89
34:BM:33:LEU:HD22	34:BM:124:LEU:HG	1.54	0.89
1:CA:699:C:C2'	1:CA:700:G:H5''	2.02	0.89
23:DB:2156:G:H2'	23:DB:2157:G:H4'	1.53	0.89
23:DB:2839:G:H4'	35:DN:49:GLU:HG2	1.55	0.89
50:B3:15:LYS:HG2	50:B3:19:GLY:HA2	1.55	0.89
5:CF:3:HIS:HA	5:CF:65:GLU:HG3	1.54	0.89
13:CN:26:LEU:HG	13:CN:44:VAL:HG22	1.55	0.89
23:DB:1639:C:H2'	23:DB:1640:A:H5''	1.52	0.89
23:DB:919:U:H2'	23:DB:920:A:C8	2.07	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:654:A:H2'	23:BB:655:A:H5''	1.55	0.88
37:DP:27:VAL:HA	37:DP:86:LYS:HE2	1.53	0.88
23:BB:2641:G:H5''	31:BJ:78:THR:HG21	1.53	0.88
23:BB:535:G:H21	38:BQ:52:ARG:NH2	1.70	0.88
32:BK:99:ILE:HG12	32:BK:115:ILE:HG13	1.52	0.88
33:BL:118:THR:HG23	33:BL:119:PRO:HD3	1.53	0.88
43:BW:48:ALA:HA	43:BW:54:ARG:H	1.38	0.88
11:CL:23:LEU:HD22	11:CL:58:ASN:HD22	1.38	0.88
1:AA:699:C:C2'	1:AA:700:G:H5''	2.02	0.88
30:BH:21:VAL:HG22	30:BH:22:LYS:H	1.38	0.88
42:BU:12:VAL:HA	42:BU:68:ASN:HD21	1.38	0.88
23:DB:2751:G:H5'	29:DG:3:VAL:HG21	1.53	0.88
34:DM:41:LEU:HD22	34:DM:95:LEU:HD13	1.53	0.88
11:AL:48:LEU:HD23	11:AL:48:LEU:H	1.34	0.88
23:BB:1021:A:H62	23:BB:1141:U:H3	1.18	0.88
26:BD:70:LYS:HA	26:BD:70:LYS:HE3	1.54	0.88
33:DL:51:GLU:HG2	33:DL:52:GLY:N	1.88	0.88
50:B3:12:ARG:NH1	50:B3:12:ARG:HB2	1.87	0.88
1:AA:74:A:C2'	1:AA:75:G:H5''	2.04	0.88
25:BC:67:LYS:HG3	25:BC:68:ARG:H	1.37	0.88
28:DF:128:SER:HB3	28:DF:154:THR:HG23	1.56	0.88
35:DN:5:LYS:HE2	35:DN:5:LYS:HA	1.55	0.88
33:BL:64:PHE:H	50:B3:24:LYS:HD2	1.36	0.88
23:BB:1141:U:H4'	23:BB:1142:A:O4'	1.74	0.88
26:BD:34:VAL:HA	26:BD:89:GLU:HB2	1.53	0.88
34:DM:29:GLY:H	34:DM:102:LEU:HD12	1.38	0.88
20:AB:67:LEU:HD12	20:AB:153:MET:HE2	1.54	0.88
33:DL:126:ARG:O	33:DL:127:VAL:HG22	1.74	0.88
44:DX:43:LEU:HB2	44:DX:45:GLN:NE2	1.89	0.88
33:BL:79:LEU:HB3	33:BL:114:GLY:H	1.35	0.88
35:BN:76:VAL:HA	35:BN:79:LEU:HD12	1.56	0.88
23:DB:929:U:H4'	45:DY:2:LYS:HE3	1.54	0.88
37:DP:32:VAL:HA	37:DP:42:PHE:HB3	1.55	0.88
23:BB:396:G:H5'	46:BZ:9:TYR:HB2	1.54	0.88
26:BD:5:VAL:HG11	26:BD:28:GLU:HA	1.54	0.88
30:BH:4:ILE:H	30:BH:37:VAL:HB	1.39	0.88
34:BM:38:ARG:NH1	34:BM:38:ARG:HB2	1.89	0.88
39:BR:37:GLU:HG2	39:BR:62:GLU:H	1.38	0.88
45:BY:15:ARG:HB3	45:BY:18:LYS:HE3	1.56	0.88
13:CN:30:ILE:HG21	13:CN:43:ALA:HB3	1.55	0.88
23:DB:1141:U:H4'	23:DB:1142:A:O4'	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1409:C:N4	1:AA:1491:G:H1	1.72	0.87
32:DK:108:ARG:HA	32:DK:116:ILE:HD13	1.56	0.87
37:BP:47:ILE:HG22	37:BP:48:ALA:H	1.37	0.87
43:BW:28:GLU:H	43:BW:61:LYS:HB2	1.37	0.87
23:DB:654:A:H2'	23:DB:655:A:H5''	1.54	0.87
39:DR:39:LEU:H	39:DR:61:ALA:HB1	1.39	0.87
39:DR:89:HIS:O	39:DR:90:ARG:HG3	1.74	0.87
38:BQ:91:ARG:H	39:BR:10:LYS:NZ	1.73	0.87
41:BT:5:GLU:HG2	41:BT:6:ARG:H	1.40	0.87
27:DE:49:ARG:HG3	27:DE:52:VAL:HG22	1.56	0.87
5:AF:38:ARG:HB3	5:AF:63:ASN:HB2	1.55	0.87
33:DL:124:GLY:H	33:DL:142:ILE:HA	1.37	0.87
2:AC:146:LYS:HD3	2:AC:204:GLY:HA2	1.56	0.87
23:BB:2147:A:H5'	23:BB:2148:G:H5'	1.55	0.87
42:BU:84:PHE:HB3	42:BU:92:VAL:HG22	1.57	0.87
23:DB:2032:G:N3	26:DD:150:GLN:HG2	1.90	0.87
26:DD:69:ALA:HB1	26:DD:92:VAL:HG13	1.56	0.87
29:BG:29:ASN:HB2	29:BG:78:VAL:HA	1.57	0.87
43:BW:19:ARG:HH22	43:BW:71:LYS:HD3	1.38	0.87
1:CA:1306:A:N6	1:CA:1331:G:H1'	1.89	0.87
26:DD:96:ILE:HG22	26:DD:98:VAL:H	1.35	0.87
27:DE:116:ASP:HB3	27:DE:185:LYS:HA	1.56	0.87
4:CE:19:ARG:HG2	4:CE:20:VAL:H	1.39	0.87
23:DB:1199:U:H2'	23:DB:1200:C:C6	2.09	0.87
40:BS:34:ASP:H	40:BS:37:THR:HG22	1.40	0.87
42:BU:40:LEU:HD13	42:BU:41:VAL:N	1.88	0.87
1:CA:1250:A:H4'	8:CI:69:GLY:H	1.38	0.87
23:BB:2354:C:H4'	43:BW:30:VAL:HG22	1.57	0.87
23:BB:2052:A:H4'	26:BD:148:GLN:N	1.90	0.87
42:BU:23:LYS:HG3	42:BU:24:VAL:HG23	1.55	0.87
23:DB:1813:G:N3	25:DC:50:THR:HG21	1.89	0.87
31:DJ:64:VAL:HG12	31:DJ:65:THR:H	1.37	0.87
37:DP:76:HIS:CD2	37:DP:76:HIS:H	1.89	0.87
43:DW:47:GLY:HA2	43:DW:71:LYS:HB3	1.57	0.87
1:AA:1323:G:H2'	1:AA:1324:A:C8	2.09	0.86
23:BB:547:A:H3'	23:BB:548:G:H8	1.40	0.86
21:CU:34:ARG:HG2	21:CU:35:GLU:H	1.39	0.86
7:AH:113:ARG:HH21	7:AH:114:ALA:HA	1.39	0.86
21:AU:36:PHE:HB3	21:AU:40:PRO:HD3	1.57	0.86
31:BJ:100:VAL:HG13	31:BJ:101:ILE:HG22	1.53	0.86
5:AF:64:VAL:HG12	5:AF:65:GLU:H	1.40	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:D4:26:ILE:HG23	51:D4:27:CYS:H	1.40	0.86
26:DD:152:PRO:HB2	26:DD:154:LYS:HE2	1.58	0.86
42:BU:43:LYS:HA	42:BU:57:ILE:HA	1.57	0.86
3:CD:25:ARG:HB2	3:CD:25:ARG:HH11	1.40	0.86
10:CK:88:PRO:HD3	21:CU:28:LEU:HD13	1.56	0.86
45:DY:2:LYS:HB2	45:DY:37:ARG:HB2	1.56	0.86
8:CI:51:LEU:HB3	8:CI:56:MET:HG2	1.58	0.86
48:B1:46:VAL:HG13	48:B1:47:ILE:HG12	1.54	0.86
23:BB:479:A:O2'	23:BB:481:G:H5'	1.75	0.86
25:BC:124:LYS:HB2	25:BC:125:PRO:CD	2.05	0.86
1:AA:1422:G:H5'	32:BK:48:PRO:HB3	1.58	0.86
3:CD:63:ILE:HG23	3:CD:64:TYR:HD1	1.40	0.86
18:CS:10:ILE:HG22	18:CS:37:SER:HB2	1.57	0.86
23:DB:1099:G:O4'	52:DI:3:LYS:C	2.14	0.86
2:AC:141:MET:HE1	2:AC:147:GLY:H	1.39	0.86
6:AG:21:LEU:H	6:AG:21:LEU:HD23	1.41	0.86
23:BB:703:U:H2'	23:BB:704:G:O4'	1.76	0.86
26:BD:130:GLN:NE2	26:BD:138:LEU:HD12	1.91	0.86
11:CL:43:LYS:HB3	11:CL:44:PRO:CD	2.06	0.86
36:BO:49:VAL:HG22	36:BO:50:ALA:H	1.38	0.86
23:DB:1098:A:H2'	52:DI:4:VAL:N	1.90	0.86
38:DQ:97:ILE:HD12	39:DR:13:ARG:HD2	1.58	0.86
7:CH:105:THR:HA	7:CH:122:GLY:HA3	1.57	0.86
33:DL:115:GLU:O	33:DL:116:VAL:HG22	1.76	0.86
42:DU:60:LYS:HE3	42:DU:61:GLU:N	1.91	0.86
11:AL:35:ARG:NH2	11:AL:75:GLU:HB3	1.89	0.86
15:AP:51:ARG:HB3	15:AP:51:ARG:NH1	1.91	0.86
31:BJ:3:THR:HG21	38:BQ:93:ILE:HG13	1.56	0.86
46:BZ:8:LYS:HD2	46:BZ:9:TYR:N	1.91	0.86
20:CB:104:LYS:HB2	20:CB:104:LYS:HZ2	1.41	0.86
25:BC:19:VAL:HG12	25:BC:20:ASN:H	1.38	0.85
39:BR:41:ILE:O	39:BR:54:VAL:HG13	1.74	0.85
25:DC:51:ARG:HH12	25:DC:54:GLY:HA3	1.40	0.85
23:BB:135:U:H2'	23:BB:136:G:C8	2.11	0.85
46:BZ:13:THR:HG23	46:BZ:25:ARG:HB3	1.56	0.85
9:CJ:6:ILE:HD11	9:CJ:79:PRO:HB3	1.58	0.85
11:CL:98:ARG:HG3	11:CL:105:GLY:HA2	1.56	0.85
33:DL:118:THR:O	33:DL:120:VAL:HG23	1.75	0.85
25:BC:74:PRO:HB2	25:BC:96:LYS:HG3	1.57	0.85
26:BD:22:ILE:N	26:BD:23:PRO:HD3	1.92	0.85
35:BN:44:LEU:HD23	35:BN:113:ILE:HG13	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CB:116:LEU:HD22	20:CB:140:LEU:HD21	1.59	0.85
25:DC:220:ARG:CZ	25:DC:220:ARG:HA	2.07	0.85
26:DD:31:ALA:HA	26:DD:51:THR:HA	1.58	0.85
52:DI:27:LEU:HD23	52:DI:27:LEU:H	1.37	0.85
43:DW:38:ARG:NH2	43:DW:40:ARG:HD3	1.91	0.85
23:BB:2052:A:C4'	26:BD:148:GLN:H	1.88	0.85
4:CE:89:THR:HG22	4:CE:91:SER:H	1.41	0.85
19:CT:68:LYS:HA	19:CT:68:LYS:HE2	1.57	0.85
23:DB:1149:G:H2'	23:DB:1150:C:C6	2.12	0.85
26:DD:204:LYS:HB3	26:DD:205:PRO:HD2	1.57	0.85
32:DK:70:ARG:HD3	32:DK:76:VAL:HG22	1.59	0.85
37:DP:38:ARG:HG3	37:DP:39:LEU:H	1.39	0.85
33:BL:29:LYS:HD2	33:BL:30:THR:H	1.42	0.85
1:CA:1296:C:H4'	1:CA:1302:C:N4	1.91	0.85
20:CB:101:THR:HA	20:CB:178:LEU:HD21	1.58	0.85
15:CP:28:ARG:HD3	15:CP:29:ASN:H	1.41	0.85
26:DD:32:ASN:HB3	26:DD:91:THR:HA	1.58	0.85
34:DM:62:LYS:H	34:DM:104:GLU:HB3	1.41	0.85
1:AA:664:G:H22	1:AA:741:G:H1	1.23	0.85
23:DB:2502:G:H5'	23:DB:2503:A:H5''	1.57	0.85
37:DP:25:VAL:HG22	37:DP:89:GLY:H	1.40	0.85
20:AB:163:ILE:HG23	20:AB:164:ASP:N	1.92	0.85
1:CA:1008:U:H5''	13:CN:23:ARG:HH21	1.41	0.85
9:AJ:35:GLN:HB2	9:AJ:78:GLU:HB2	1.57	0.85
34:BM:7:THR:HG22	34:BM:8:LYS:H	1.41	0.85
1:CA:1306:A:H61	1:CA:1331:G:H1'	1.40	0.85
2:CC:69:THR:HG22	2:CC:71:ARG:H	1.38	0.85
33:DL:39:LYS:HA	33:DL:39:LYS:HZ2	1.40	0.85
19:AT:30:PHE:HB3	19:AT:53:MET:HB3	1.59	0.85
28:DF:107:VAL:HG12	28:DF:108:PRO:HD3	1.57	0.85
6:AG:112:ASP:HB2	6:AG:118:ARG:HG2	1.58	0.85
31:BJ:101:ILE:HG23	31:BJ:102:GLU:H	1.42	0.85
14:CO:69:LEU:HD11	14:CO:76:ARG:HB3	1.57	0.85
27:DE:3:LEU:HD22	27:DE:119:ILE:HD11	1.57	0.85
31:DJ:35:ARG:NH1	31:DJ:40:HIS:H	1.75	0.85
33:BL:77:ILE:HG13	33:BL:108:ALA:HB1	1.59	0.84
26:DD:14:ILE:HD12	37:DP:78:PRO:HG2	1.59	0.84
39:DR:47:VAL:HG22	39:DR:48:LYS:H	1.40	0.84
43:DW:67:LYS:HG2	43:DW:71:LYS:HB2	1.59	0.84
6:AG:97:ALA:HA	6:AG:100:MET:HE3	1.59	0.84
29:BG:15:ASP:HB3	29:BG:26:LYS:HB2	1.58	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:664:G:H22	1:CA:741:G:H1	1.23	0.84
1:CA:9:G:H5'	4:CE:107:GLY:HA3	1.55	0.84
3:CD:28:ASP:HB3	3:CD:33:ILE:HD12	1.58	0.84
10:AK:126:ARG:HB2	21:AU:33:ARG:HD2	1.59	0.84
34:BM:70:ASP:H	34:BM:71:LYS:HZ2	1.23	0.84
1:CA:840:C:H2'	1:CA:842:U:H5''	1.60	0.84
1:AA:1101:A:H4'	1:AA:1102:A:O5'	1.75	0.84
12:AM:33:LEU:HD22	12:AM:38:ILE:HB	1.59	0.84
23:BB:662:G:H4'	33:BL:24:GLY:N	1.92	0.84
36:BO:16:ARG:HE	36:BO:16:ARG:HA	1.42	0.84
42:BU:3:LYS:NZ	42:BU:3:LYS:HB2	1.90	0.84
20:CB:16:GLY:HA3	20:CB:39:ILE:HA	1.59	0.84
15:CP:67:ILE:HG13	15:CP:71:VAL:HG13	1.59	0.84
26:DD:122:VAL:HA	26:DD:128:ARG:HG3	1.59	0.84
1:AA:74:A:H2'	1:AA:75:G:H5''	1.58	0.84
23:BB:2306:C:H42	28:BF:38:GLY:H	1.24	0.84
23:BB:2502:G:H5'	23:BB:2503:A:H5''	1.57	0.84
26:BD:38:LYS:HG2	26:BD:39:ASP:H	1.42	0.84
11:CL:68:GLY:HA3	11:CL:106:VAL:HG21	1.59	0.84
15:CP:44:SER:HB3	15:CP:46:LYS:HG2	1.60	0.84
21:CU:39:LYS:H	21:CU:40:PRO:HD2	1.43	0.84
23:DB:443:A:OP1	27:DE:40:ARG:HA	1.76	0.84
1:AA:840:C:H2'	1:AA:842:U:H5''	1.59	0.84
6:AG:3:ARG:NH1	6:AG:3:ARG:HB3	1.92	0.84
27:BE:128:ALA:CB	27:BE:129:PRO:HD3	2.07	0.84
1:CA:781:A:H2'	1:CA:782:A:H5'	1.60	0.84
23:DB:1021:A:H61	23:DB:1142:A:N6	1.73	0.84
12:AM:15:VAL:HG23	12:AM:33:LEU:HD12	1.60	0.84
44:DX:18:LEU:HA	44:DX:22:LEU:HD12	1.58	0.84
46:DZ:24:ILE:HD13	46:DZ:24:ILE:H	1.41	0.84
14:AO:25:GLU:HG3	14:AO:80:LEU:HD12	1.59	0.84
23:BB:2619:C:H5'	26:BD:157:LYS:HE2	1.59	0.84
35:BN:101:GLY:HA2	35:BN:109:PRO:HA	1.58	0.84
37:BP:59:THR:HG23	37:BP:76:HIS:CD2	2.12	0.84
27:DE:47:LYS:HE3	27:DE:52:VAL:HG23	1.60	0.84
30:DH:37:VAL:H	30:DH:38:PRO:HD2	1.43	0.84
31:DJ:135:GLN:HE22	31:DJ:137:PRO:HB2	1.43	0.84
10:AK:86:LYS:HB2	10:AK:112:VAL:HG23	1.59	0.84
38:BQ:30:VAL:HG12	38:BQ:32:ARG:H	1.43	0.84
38:BQ:92:LYS:CG	38:BQ:93:ILE:H	1.90	0.84
11:CL:84:GLY:H	11:CL:94:TYR:HA	1.42	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:84:ASN:HD22	3:AD:87:GLU:H	1.26	0.84
25:BC:155:ARG:HE	25:BC:157:ALA:HB3	1.41	0.84
23:BB:2024:G:H4'	26:BD:154:LYS:HD3	1.60	0.84
28:BF:66:ILE:HG22	28:BF:83:PRO:HB2	1.59	0.84
33:BL:116:VAL:HG12	33:BL:117:THR:H	1.42	0.84
23:DB:1799:G:N7	25:DC:178:GLY:HA3	1.92	0.84
1:CA:1422:G:H5''	32:DK:48:PRO:HB3	1.57	0.84
1:AA:1330:U:H2'	1:AA:1331:G:H5'	1.59	0.83
1:AA:33:A:H1'	11:AL:27:PRO:HG3	1.59	0.83
30:BH:130:VAL:HB	30:BH:142:VAL:HB	1.60	0.83
24:BV:80:HIS:CD2	24:BV:83:LYS:H	1.94	0.83
1:CA:108:G:O6	19:CT:9:ARG:HG2	1.78	0.83
2:CC:179:ALA:HA	2:CC:205:GLU:HA	1.59	0.83
30:DH:10:ALA:O	30:DH:11:ASN:HB3	1.78	0.83
12:AM:3:ILE:HA	12:AM:56:ARG:HG2	1.60	0.83
16:AQ:67:SER:HB2	16:AQ:70:LYS:HB3	1.60	0.83
35:BN:19:ALA:HA	35:BN:22:ARG:HG2	1.58	0.83
8:CI:54:VAL:HG11	8:CI:86:LEU:HD13	1.58	0.83
10:CK:113:THR:HG21	21:CU:28:LEU:HD11	1.59	0.83
20:AB:33:ALA:HA	20:AB:38:HIS:HA	1.60	0.83
25:BC:136:VAL:HG12	25:BC:165:ALA:HA	1.60	0.83
20:CB:113:LEU:HD12	20:CB:143:LEU:HD22	1.60	0.83
23:DB:2314:A:H1'	28:DF:154:THR:HG21	1.60	0.83
28:DF:70:ARG:HA	28:DF:80:GLN:HE21	1.42	0.83
30:DH:31:VAL:HB	30:DH:32:PRO:HD3	1.60	0.83
37:DP:52:ARG:HB3	37:DP:60:VAL:HG11	1.60	0.83
13:AN:30:ILE:HD12	13:AN:30:ILE:H	1.43	0.83
23:BB:775:G:H4'	23:BB:776:G:H5'	1.59	0.83
32:BK:108:ARG:NH2	37:BP:36:LYS:HG3	1.92	0.83
12:CM:38:ILE:HG13	12:CM:55:LEU:HD21	1.60	0.83
23:DB:996:A:H5''	38:DQ:93:ILE:HG21	1.58	0.83
34:DM:5:LYS:HE3	34:DM:6:ARG:H	1.43	0.83
38:DQ:73:ILE:HG13	38:DQ:74:SER:H	1.43	0.83
40:DS:25:ARG:HD2	40:DS:26:GLY:N	1.94	0.83
1:AA:1005:A:H2'	1:AA:1006:G:O4'	1.78	0.83
25:BC:42:ARG:HG2	25:BC:43:ASN:N	1.89	0.83
4:CE:157:GLY:H	7:CH:43:GLY:HA3	1.41	0.83
13:CN:96:LYS:HG2	13:CN:97:LYS:H	1.44	0.83
31:DJ:84:ILE:HD12	31:DJ:85:LYS:N	1.92	0.83
1:AA:781:A:H2'	1:AA:782:A:H5'	1.60	0.83
23:BB:1024:G:H3'	23:BB:1025:G:H5''	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2297:A:N6	23:BB:2319:G:H3'	1.93	0.83
44:BX:35:GLY:HA3	44:BX:40:SER:HA	1.60	0.83
39:DR:66:HIS:HA	39:DR:98:ILE:HD13	1.57	0.83
7:AH:5:PRO:HB2	7:AH:32:LYS:NZ	1.92	0.83
23:BB:445:C:H5'	27:BE:47:LYS:HZ3	1.41	0.83
20:CB:100:LEU:HB3	20:CB:178:LEU:HD11	1.61	0.83
1:AA:967:C:H3'	1:AA:968:A:H5'	1.59	0.83
3:AD:160:LEU:H	3:AD:160:LEU:HD13	1.44	0.83
23:BB:1060:U:C2	23:BB:1088:A:N7	2.47	0.83
33:BL:84:LYS:HA	33:BL:84:LYS:NZ	1.93	0.83
4:CE:36:THR:HG21	4:CE:63:MET:HG2	1.61	0.83
23:DB:162:U:O2'	23:DB:163:C:H5'	1.79	0.83
23:DB:2143:C:H2'	23:DB:2144:G:O4'	1.77	0.83
23:DB:2309:A:H3'	23:DB:2310:C:H5''	1.60	0.83
40:DS:23:LEU:HD13	40:DS:25:ARG:HH22	1.44	0.83
42:DU:9:GLU:HB2	42:DU:71:ILE:HB	1.60	0.83
1:AA:1320:C:H42	18:AS:35:ARG:HD3	1.43	0.83
29:BG:27:GLY:HA3	29:BG:78:VAL:HB	1.60	0.83
34:BM:82:MET:HB2	34:BM:84:LYS:NZ	1.94	0.83
15:CP:3:THR:HG22	15:CP:66:THR:HB	1.61	0.83
18:CS:32:THR:HG22	18:CS:33:TRP:H	1.43	0.83
23:DB:775:G:H4'	23:DB:776:G:H5'	1.59	0.83
23:BB:2845:U:O3'	37:BP:54:LEU:HG	1.78	0.83
40:BS:90:LYS:HZ2	40:BS:90:LYS:N	1.76	0.83
44:BX:14:LEU:HD21	44:BX:57:LEU:HB2	1.61	0.83
19:CT:61:ALA:HA	19:CT:67:HIS:H	1.44	0.83
41:DT:47:VAL:HG22	41:DT:53:VAL:HG21	1.61	0.83
23:BB:1913:A:H1'	23:BB:1914:C:OP1	1.79	0.82
44:BX:38:GLN:H	44:BX:38:GLN:NE2	1.75	0.82
7:CH:11:THR:HG22	7:CH:14:ARG:HH12	1.44	0.82
23:DB:1060:U:C2	23:DB:1088:A:N7	2.47	0.82
23:DB:136:G:C2	41:DT:3:ARG:CZ	2.62	0.82
9:AJ:8:ILE:HD12	9:AJ:100:ILE:HG22	1.61	0.82
23:BB:37:C:H2'	27:BE:46:GLN:CD	1.99	0.82
18:CS:15:LEU:HA	18:CS:18:VAL:HG12	1.60	0.82
23:DB:90:U:H3'	23:DB:91:A:H5''	1.59	0.82
35:DN:41:ALA:HB1	35:DN:113:ILE:HD11	1.61	0.82
23:BB:630:G:H1	33:BL:69:ARG:NH2	1.77	0.82
25:BC:138:SER:HA	25:BC:162:GLN:HG2	1.61	0.82
45:BY:9:THR:HB	45:BY:54:VAL:HA	1.58	0.82
22:DA:30:C:H2'	22:DA:31:C:H5'	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DN:37:THR:HG22	35:DN:39:PRO:HD2	1.59	0.82
35:DN:85:PRO:HA	35:DN:88:ALA:HB2	1.62	0.82
23:BB:997:G:OP1	38:BQ:92:LYS:HD2	1.79	0.82
45:BY:6:ILE:HG22	45:BY:56:VAL:HG23	1.62	0.82
5:CF:99:ALA:HA	17:CR:23:LYS:HE2	1.62	0.82
28:DF:64:PRO:HA	28:DF:88:VAL:HG22	1.61	0.82
23:DB:7:G:H5'	31:DJ:134:ALA:O	1.80	0.82
43:DW:3:LYS:HA	43:DW:3:LYS:NZ	1.94	0.82
2:AC:78:LYS:HG2	2:AC:81:GLU:HB2	1.58	0.82
12:AM:64:VAL:HA	12:AM:68:LEU:HD12	1.61	0.82
23:DB:2574:G:H21	26:DD:147:GLY:CA	1.92	0.82
23:DB:742:A:H2'	23:DB:743:A:H8	1.43	0.82
23:DB:1060:U:H5	52:DI:131:THR:HG22	1.43	0.82
27:BE:88:ARG:HH11	27:BE:88:ARG:HB2	1.45	0.82
27:DE:142:ALA:H	27:DE:185:LYS:NZ	1.78	0.82
36:DO:56:LYS:HE2	36:DO:81:ARG:NE	1.94	0.82
23:DB:2334:U:H3	43:DW:74:LYS:HE2	1.44	0.82
23:BB:1639:C:C2'	23:BB:1640:A:H5''	2.09	0.82
23:BB:742:A:H2'	23:BB:743:A:H8	1.43	0.82
43:BW:27:GLY:HA2	43:BW:60:ALA:HA	1.60	0.82
52:DI:72:THR:HG21	52:DI:112:LYS:HA	1.62	0.82
42:DU:13:LEU:H	42:DU:13:LEU:HD12	1.45	0.82
23:BB:287:G:H2'	23:BB:288:U:C6	2.15	0.82
30:BH:128:HIS:HB3	30:BH:144:VAL:HB	1.61	0.82
52:BI:55:PRO:HD3	52:BI:74:PRO:HD3	1.60	0.82
23:DB:2336:A:H1'	23:DB:2337:G:OP1	1.79	0.82
10:AK:83:VAL:HG23	10:AK:109:ILE:HA	1.60	0.82
17:AR:52:ARG:HH11	17:AR:52:ARG:HB3	1.45	0.82
33:BL:29:LYS:HD2	33:BL:31:GLY:N	1.95	0.82
42:BU:25:LYS:HG3	42:BU:35:VAL:HA	1.62	0.82
42:BU:7:ASP:H	42:BU:24:VAL:HA	1.45	0.82
6:CG:129:ASN:HB3	6:CG:134:VAL:HG21	1.61	0.82
10:CK:16:SER:HA	10:CK:78:ILE:HA	1.61	0.82
52:DI:105:LEU:HD13	52:DI:129:GLU:HG2	1.61	0.82
15:AP:12:LYS:HD2	15:AP:13:LYS:HE2	1.61	0.82
23:BB:2297:A:H61	23:BB:2319:G:C3'	1.93	0.82
33:BL:109:LYS:HA	33:BL:126:ARG:HA	1.62	0.82
45:BY:7:THR:HA	45:BY:34:THR:HA	1.61	0.82
30:DH:114:GLU:HB2	30:DH:133:GLN:HG3	1.60	0.82
21:AU:20:ARG:HA	21:AU:24:LYS:HG3	1.59	0.81
23:BB:2336:A:H1'	23:BB:2337:G:OP1	1.80	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1639:C:C2'	23:DB:1640:A:H5''	2.09	0.81
23:DB:2633:G:H1'	26:DD:62:LYS:HG3	1.61	0.81
28:DF:96:TRP:O	28:DF:100:GLU:HG3	1.78	0.81
18:AS:4:LEU:HD13	18:AS:9:PHE:H	1.45	0.81
32:BK:35:VAL:HG23	32:BK:36:GLY:H	1.44	0.81
36:BO:17:LYS:HE2	43:BW:77:LYS:HD2	1.61	0.81
30:DH:122:LEU:HD13	30:DH:146:VAL:HG22	1.60	0.81
42:DU:91:LYS:HD3	42:DU:93:ARG:HE	1.45	0.81
29:DG:171:LYS:HD3	29:DG:172:GLU:O	1.79	0.81
33:DL:103:ILE:HG13	33:DL:104:GLN:H	1.45	0.81
37:DP:28:LYS:HD3	37:DP:44:GLY:HA3	1.63	0.81
8:AI:112:ARG:HB2	8:AI:112:ARG:NH1	1.95	0.81
8:AI:40:ARG:H	8:AI:44:ARG:HE	1.26	0.81
23:BB:340:A:H2'	23:BB:341:C:O4'	1.81	0.81
30:BH:66:ASN:HA	30:BH:138:VAL:HG21	1.61	0.81
36:BO:64:TYR:HE1	36:BO:74:VAL:HG11	1.46	0.81
36:BO:87:ILE:HG13	36:BO:88:LYS:H	1.43	0.81
23:DB:857:G:C2'	23:DB:858:G:H5'	2.10	0.81
23:DB:1064:C:H4'	52:DI:90:GLY:HA2	1.60	0.81
3:AD:36:ALA:HA	3:AD:41:GLY:HA3	1.62	0.81
23:BB:279:A:N6	23:BB:361:G:H1'	1.96	0.81
28:BF:126:ASN:HD22	28:BF:126:ASN:N	1.78	0.81
23:BB:200:U:H5''	46:BZ:20:ASN:HB3	1.62	0.81
23:DB:899:A:H3'	23:DB:900:A:H8	1.45	0.81
31:DJ:25:LEU:HD11	31:DJ:63:ALA:H	1.45	0.81
39:DR:22:LEU:HD12	39:DR:24:LYS:H	1.43	0.81
36:BO:8:ILE:HG22	36:BO:10:ARG:HG2	1.63	0.81
42:BU:28:LEU:HB2	42:BU:31:GLY:HA3	1.61	0.81
1:CA:1086:U:H3	1:CA:1099:G:H22	1.25	0.81
20:CB:163:ILE:HG23	20:CB:164:ASP:H	1.46	0.81
38:DQ:87:VAL:HB	39:DR:54:VAL:HG21	1.62	0.81
43:DW:60:ALA:HB3	43:DW:80:SER:HA	1.63	0.81
25:BC:163:ILE:HG22	25:BC:164:VAL:H	1.44	0.81
31:BJ:100:VAL:HG22	31:BJ:101:ILE:H	1.45	0.81
37:BP:92:ARG:NH2	37:BP:110:LYS:HE2	1.95	0.81
37:BP:47:ILE:HG22	37:BP:49:ILE:H	1.45	0.81
40:BS:71:VAL:HA	40:BS:106:VAL:O	1.80	0.81
43:BW:42:THR:HB	43:BW:65:LYS:HA	1.60	0.81
25:DC:172:THR:O	25:DC:173:LEU:HB2	1.81	0.81
3:AD:194:ILE:HD13	3:AD:195:ASN:N	1.96	0.81
11:AL:56:LEU:HD11	11:AL:81:ILE:HD12	1.61	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1177:G:H2'	23:BB:1178:C:C6	2.16	0.81
38:BQ:7:VAL:HG23	38:BQ:8:ILE:H	1.43	0.81
23:DB:2511:U:H5''	26:DD:129:THR:HG23	1.62	0.81
23:DB:589:U:H4'	27:DE:87:ALA:HB2	1.61	0.81
12:AM:11:HIS:H	12:AM:44:ILE:HD11	1.44	0.81
16:AQ:27:PHE:HB3	16:AQ:36:PHE:HB3	1.63	0.81
46:BZ:26:SER:O	46:BZ:28:VAL:HG13	1.81	0.81
3:CD:162:GLU:HG3	3:CD:163:GLN:HG3	1.63	0.81
8:CI:45:MET:O	8:CI:49:GLN:HG3	1.80	0.81
23:DB:250:G:H5'	50:D3:7:ARG:HG2	1.61	0.81
52:DI:105:LEU:HD11	52:DI:139:VAL:HG21	1.62	0.81
1:AA:238:A:H2'	1:AA:239:U:H5''	1.61	0.81
38:BQ:73:ILE:HG23	38:BQ:77:LYS:HB2	1.62	0.81
19:CT:45:ALA:HA	19:CT:48:LYS:HB3	1.63	0.81
23:DB:458:G:N2	23:DB:469:G:H2'	1.95	0.81
23:DB:717:C:H3'	23:DB:718:A:H5''	1.61	0.81
25:DC:140:VAL:HA	25:DC:191:LEU:HA	1.62	0.81
18:AS:5:LYS:N	18:AS:5:LYS:HE3	1.96	0.81
31:BJ:16:TYR:HB2	31:BJ:54:ILE:HG22	1.63	0.81
32:BK:64:ARG:H	32:BK:83:ALA:HB3	1.44	0.81
50:D3:7:ARG:NH1	50:D3:7:ARG:HA	1.95	0.81
27:DE:6:LYS:HD3	27:DE:7:ASP:H	1.46	0.81
16:AQ:20:ILE:HG13	16:AQ:45:VAL:HB	1.61	0.80
37:BP:25:VAL:HG13	37:BP:87:ARG:HA	1.61	0.80
1:CA:1238:A:H5'	1:CA:1336:C:H41	1.45	0.80
41:DT:34:VAL:HG21	41:DT:43:ILE:HD12	1.62	0.80
46:DZ:59:ARG:HB3	46:DZ:63:ARG:HB2	1.61	0.80
1:AA:1323:G:H2'	1:AA:1324:A:H8	1.44	0.80
35:BN:96:ARG:HH21	35:BN:118:ARG:HB3	1.46	0.80
23:BB:580:U:O3'	38:BQ:30:VAL:HG13	1.80	0.80
49:D2:35:ARG:HH22	49:D2:44:VAL:HG22	1.46	0.80
23:DB:2377:A:N6	36:DO:13:ARG:HH21	1.78	0.80
2:AC:87:ARG:HG3	2:AC:100:ILE:HG22	1.63	0.80
23:BB:265:A:O2'	23:BB:266:G:H4'	1.81	0.80
25:BC:22:GLU:HA	25:BC:202:ARG:CZ	2.11	0.80
26:BD:120:GLY:H	26:BD:123:LYS:HG3	1.43	0.80
36:BO:1:MET:HG2	36:BO:3:LYS:H	1.45	0.80
38:BQ:16:ILE:HD12	38:BQ:17:LEU:N	1.96	0.80
39:BR:75:VAL:H	39:BR:90:ARG:NE	1.79	0.80
47:D0:27:LEU:H	47:D0:27:LEU:HD22	1.45	0.80
33:DL:124:GLY:N	33:DL:142:ILE:HA	1.95	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DQ:87:VAL:HG12	38:DQ:89:ILE:HD13	1.62	0.80
23:BB:704:G:H2'	23:BB:726:G:H22	1.45	0.80
25:BC:194:VAL:HG12	25:BC:195:GLY:H	1.45	0.80
23:BB:1568:G:OP1	25:BC:61:TYR:HB2	1.82	0.80
29:BG:153:PRO:HG2	29:BG:162:ARG:HB3	1.63	0.80
1:CA:238:A:H2'	1:CA:239:U:H5''	1.61	0.80
23:DB:354:A:H2'	23:DB:355:U:C6	2.17	0.80
30:DH:115:VAL:HG22	30:DH:117:LEU:H	1.45	0.80
32:DK:2:ILE:HD13	32:DK:6:THR:HG21	1.61	0.80
33:DL:7:SER:CB	33:DL:8:PRO:HD3	2.07	0.80
37:DP:50:ARG:NH1	37:DP:50:ARG:HB2	1.97	0.80
3:AD:84:ASN:ND2	3:AD:87:GLU:H	1.79	0.80
8:AI:35:GLU:HA	8:AI:39:GLY:HA3	1.64	0.80
23:BB:1178:C:H2'	23:BB:1179:G:C8	2.16	0.80
31:BJ:35:ARG:HE	31:BJ:39:LYS:HB3	1.46	0.80
31:BJ:71:ASP:CG	31:BJ:72:LYS:H	1.85	0.80
29:DG:40:VAL:HG22	29:DG:51:PHE:HE2	1.44	0.80
39:DR:64:VAL:HG22	39:DR:65:ALA:H	1.46	0.80
2:AC:176:THR:HB	2:AC:179:ALA:HB2	1.63	0.80
3:AD:94:GLU:HG3	3:AD:103:ARG:HH12	1.46	0.80
43:BW:36:ILE:HD12	43:BW:69:GLU:OE2	1.82	0.80
1:CA:1451:U:H5''	1:CA:1452:C:OP2	1.81	0.80
32:DK:78:ARG:HH22	37:DP:62:LYS:NZ	1.80	0.80
1:AA:1201:A:H1'	1:AA:1202:U:OP2	1.81	0.80
23:DB:1007:C:H5''	31:DJ:37:ARG:HH12	1.47	0.80
37:DP:69:VAL:HG13	37:DP:70:GLU:H	1.45	0.80
1:AA:1026:G:H2'	1:AA:1027:C:H6	1.47	0.80
23:BB:144:A:H2'	23:BB:145:C:H6	1.47	0.80
31:BJ:25:LEU:HD12	31:BJ:62:VAL:HA	1.63	0.80
39:BR:4:VAL:HA	39:BR:12:HIS:HB3	1.63	0.80
41:BT:12:ARG:HH21	44:BX:29:ARG:HH12	1.27	0.80
6:CG:112:ASP:HB2	6:CG:118:ARG:HG2	1.62	0.80
26:DD:37:VAL:HG13	26:DD:42:ASN:HB3	1.63	0.80
28:DF:98:PHE:HA	28:DF:101:ARG:HE	1.45	0.80
31:DJ:121:LYS:N	31:DJ:121:LYS:HE3	1.97	0.80
33:DL:135:ILE:HG22	33:DL:138:ALA:HB3	1.63	0.80
1:AA:946:A:H2'	1:AA:947:G:C8	2.16	0.80
52:BI:106:GLN:O	52:BI:110:GLN:HG3	1.81	0.80
39:BR:6:GLN:HA	39:BR:10:LYS:HA	1.64	0.80
39:BR:58:VAL:HG13	39:BR:59:ILE:HG12	1.61	0.80
4:CE:36:THR:HG22	4:CE:37:VAL:H	1.46	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:120:VAL:HG12	27:DE:121:VAL:H	1.45	0.80
30:DH:31:VAL:HA	30:DH:36:ALA:HA	1.63	0.80
31:DJ:25:LEU:HG	31:DJ:64:VAL:H	1.47	0.80
1:AA:1326:U:H2'	1:AA:1327:C:H6	1.47	0.80
25:BC:88:ALA:HB1	25:BC:156:SER:HB3	1.64	0.80
32:BK:88:ASN:ND2	32:BK:91:SER:HB2	1.97	0.80
33:BL:60:ARG:HA	50:B3:11:LYS:HE2	1.63	0.80
40:BS:50:VAL:HG11	40:BS:103:ILE:HG21	1.62	0.80
23:BB:100:U:H4'	42:BU:90:LYS:CE	2.12	0.80
10:CK:70:ALA:HB1	10:CK:74:LYS:HD2	1.64	0.80
21:CU:14:ALA:HB3	21:CU:16:ARG:NH1	1.97	0.80
23:DB:1099:G:H5''	52:DI:2:LYS:C	2.01	0.80
26:DD:109:VAL:HG11	26:DD:193:VAL:HG12	1.64	0.80
33:DL:109:LYS:HG2	33:DL:126:ARG:HD3	1.63	0.80
37:DP:64:SER:HA	37:DP:71:ARG:HB2	1.62	0.80
23:BB:1131:G:H1'	23:BB:1133:A:H62	1.47	0.79
23:BB:279:A:C2	23:BB:362:A:H4'	2.17	0.79
23:BB:692:C:C5'	25:BC:42:ARG:HB2	2.12	0.79
31:BJ:25:LEU:HG	31:BJ:64:VAL:H	1.47	0.79
31:BJ:72:LYS:HB2	31:BJ:90:GLU:HB2	1.61	0.79
41:BT:1:MET:SD	41:BT:2:ILE:HG12	2.21	0.79
7:CH:55:LYS:HA	7:CH:55:LYS:HZ2	1.45	0.79
48:D1:12:SER:HA	48:D1:50:GLU:HA	1.63	0.79
23:DB:1098:A:H3'	52:DI:3:LYS:CB	2.11	0.79
38:DQ:69:ARG:HB3	38:DQ:69:ARG:HH11	1.47	0.79
45:DY:10:ARG:HA	45:DY:31:ILE:HD12	1.64	0.79
8:AI:110:VAL:HG12	8:AI:111:GLU:H	1.45	0.79
23:BB:162:U:H4'	23:BB:163:C:OP1	1.82	0.79
27:BE:153:LEU:HD13	27:BE:193:VAL:HG21	1.65	0.79
33:BL:84:LYS:HA	33:BL:84:LYS:HZ2	1.45	0.79
39:BR:41:ILE:HD12	39:BR:41:ILE:O	1.83	0.79
4:CE:33:THR:HG22	4:CE:51:LYS:HG2	1.64	0.79
52:DI:21:PRO:HB2	52:DI:22:PRO:HD3	1.64	0.79
37:DP:25:VAL:HG13	37:DP:88:ARG:H	1.46	0.79
26:BD:4:LEU:HD13	26:BD:79:LEU:HD11	1.65	0.79
30:BH:83:LYS:HB3	30:BH:91:PHE:HD1	1.43	0.79
34:BM:42:THR:HA	34:BM:92:TRP:CD1	2.16	0.79
37:BP:2:ASN:H	37:BP:2:ASN:HD22	1.30	0.79
2:CC:18:ASN:HA	2:CC:55:VAL:HG12	1.65	0.79
23:DB:85:G:H5'	42:DU:28:LEU:HB3	1.61	0.79
41:BT:12:ARG:HH21	44:BX:29:ARG:NH1	1.80	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:60:A:H2'	19:CT:4:LYS:HE3	1.64	0.79
23:DB:1176:U:H2'	23:DB:1177:G:C8	2.17	0.79
28:DF:140:ILE:HG22	28:DF:142:TYR:H	1.47	0.79
33:DL:81:ASP:HA	33:DL:84:LYS:HD3	1.64	0.79
23:BB:972:A:H3'	23:BB:973:A:H5''	1.63	0.79
41:BT:8:LEU:HA	44:BX:19:LEU:HD21	1.64	0.79
2:CC:120:THR:HG23	2:CC:188:ALA:HB2	1.63	0.79
47:D0:53:VAL:HG13	47:D0:54:ILE:HG13	1.65	0.79
23:DB:320:A:H4'	23:DB:322:A:N7	1.97	0.79
28:DF:133:GLU:N	28:DF:150:GLY:HA2	1.98	0.79
30:DH:114:GLU:HB3	30:DH:133:GLN:HE21	1.47	0.79
23:DB:1006:C:H5''	31:DJ:34:ARG:HE	1.48	0.79
23:BB:1197:G:H2'	23:BB:1198:U:C6	2.18	0.79
23:BB:1813:G:H21	25:BC:51:ARG:HG2	1.46	0.79
25:BC:50:THR:HG22	25:BC:51:ARG:HG3	1.65	0.79
1:CA:1003:G:N2	1:CA:1005:A:H5'	1.97	0.79
1:CA:1311:A:H62	18:CS:1:PRO:HD3	1.48	0.79
20:CB:202:ASN:ND2	20:CB:203:ASP:H	1.79	0.79
4:CE:158:LYS:HB3	7:CH:63:LYS:HD3	1.65	0.79
13:CN:63:CYS:HB3	13:CN:68:ARG:H	1.47	0.79
22:DA:48:U:H2'	22:DA:49:C:H6	1.44	0.79
41:DT:56:GLU:O	41:DT:57:VAL:HG22	1.82	0.79
30:BH:3:VAL:HB	30:BH:37:VAL:HB	1.64	0.79
6:CG:23:ALA:O	6:CG:26:VAL:HG22	1.82	0.79
8:CI:17:ARG:HB3	8:CI:65:THR:HB	1.65	0.79
25:DC:63:ILE:HG21	25:DC:90:ILE:HD11	1.64	0.79
27:DE:53:THR:HB	27:DE:74:LYS:HG2	1.63	0.79
33:DL:44:GLY:HA2	33:DL:47:ARG:HE	1.46	0.79
24:DV:70:ILE:HD12	24:DV:71:LYS:H	1.47	0.79
45:DY:45:GLY:HA2	45:DY:48:ASN:HD22	1.46	0.79
22:BA:49:C:H2'	22:BA:50:A:H8	1.48	0.79
23:BB:3:U:HO2'	23:BB:4:U:H6	1.29	0.79
25:BC:128:THR:HA	25:BC:190:THR:HA	1.62	0.79
25:BC:3:VAL:HG12	25:BC:4:LYS:H	1.48	0.79
26:BD:11:MET:HG2	26:BD:12:THR:H	1.45	0.79
27:BE:118:LEU:HD23	27:BE:118:LEU:H	1.45	0.79
23:BB:37:C:H2'	27:BE:46:GLN:HG2	1.63	0.79
33:BL:61:LEU:CB	33:BL:62:PRO:HD2	2.12	0.79
13:CN:50:LEU:H	13:CN:51:PRO:CD	1.95	0.79
16:CQ:66:LEU:HD12	16:CQ:66:LEU:H	1.47	0.79
28:DF:35:LEU:HB3	28:DF:151:LEU:HD11	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DI:11:GLN:HG2	52:DI:55:PRO:HB3	1.64	0.79
34:DM:4:PRO:HD2	34:DM:43:ALA:HB1	1.65	0.79
44:DX:4:LYS:O	44:DX:7:ARG:HG2	1.82	0.79
1:AA:17:U:H2'	1:AA:18:C:C6	2.18	0.79
9:AJ:40:ILE:HG13	9:AJ:73:LEU:HB3	1.65	0.79
16:AQ:10:ARG:CZ	16:AQ:11:VAL:H	1.95	0.79
52:BI:27:LEU:HD12	52:BI:32:VAL:HG11	1.64	0.79
33:BL:27:LEU:HD12	33:BL:27:LEU:H	1.47	0.79
34:BM:41:LEU:HB2	34:BM:93:VAL:HG22	1.64	0.79
2:CC:137:VAL:HA	2:CC:148:ILE:HD13	1.63	0.79
5:CF:47:LEU:HD13	5:CF:51:ILE:HG22	1.62	0.79
33:DL:4:ASN:O	33:DL:6:LEU:HD22	1.82	0.79
9:AJ:88:MET:HB2	9:AJ:89:ARG:HH12	1.47	0.79
51:B4:2:LYS:HG2	51:B4:36:ARG:HB3	1.64	0.79
27:BE:113:VAL:HG23	27:BE:117:ARG:CZ	2.13	0.79
34:BM:126:ILE:HG13	34:BM:127:LYS:H	1.47	0.79
1:CA:946:A:H2'	1:CA:947:G:C8	2.18	0.79
23:DB:1099:G:H5''	52:DI:3:LYS:N	1.98	0.79
23:DB:2800:A:H2'	23:DB:2801:G:O4'	1.83	0.79
29:DG:120:ILE:HG12	29:DG:134:GLY:HA3	1.65	0.79
30:DH:68:ARG:HE	30:DH:71:LYS:HE2	1.47	0.79
35:DN:8:ARG:HD2	35:DN:46:ARG:NE	1.98	0.79
46:DZ:66:ILE:HB	46:DZ:67:PRO:HD3	1.64	0.79
1:AA:436:C:O2'	1:AA:437:U:H5'	1.83	0.78
6:AG:4:ARG:HE	6:AG:5:VAL:H	1.30	0.78
23:BB:86:G:OP1	42:BU:29:SER:HB2	1.83	0.78
43:BW:28:GLU:HB3	43:BW:30:VAL:HG23	1.65	0.78
46:BZ:44:PHE:H	46:BZ:44:PHE:HD1	1.30	0.78
23:DB:1802:A:H4'	25:DC:255:LYS:HE2	1.63	0.78
25:BC:140:VAL:O	25:BC:193:GLU:HB2	1.82	0.78
26:BD:37:VAL:HG22	26:BD:46:ARG:HE	1.48	0.78
37:BP:108:ARG:HD2	37:BP:108:ARG:H	1.48	0.78
43:BW:66:VAL:HG13	43:BW:67:LYS:H	1.47	0.78
23:DB:2309:A:H5'	23:DB:2310:C:OP2	1.83	0.78
37:DP:30:TRP:HB2	37:DP:83:ILE:HB	1.64	0.78
13:AN:65:GLN:HB2	13:AN:78:LEU:HD22	1.63	0.78
23:BB:2216:G:H2'	23:BB:2217:G:H8	1.47	0.78
23:BB:222:A:H61	23:BB:232:G:H1'	1.45	0.78
23:BB:2511:U:H4'	26:BD:128:ARG:CZ	2.13	0.78
23:BB:2800:A:H2'	23:BB:2801:G:O4'	1.84	0.78
30:BH:68:ARG:HG3	30:BH:134:VAL:HG11	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:BL:60:ARG:HD2	50:B3:7:ARG:NE	1.98	0.78
42:BU:35:VAL:HG12	42:BU:36:GLU:H	1.47	0.78
23:DB:1080:A:H2'	23:DB:1081:U:H6	1.45	0.78
39:DR:40:MET:HG2	39:DR:54:VAL:HG13	1.65	0.78
23:DB:922:C:H1'	43:DW:22:VAL:HG21	1.66	0.78
1:AA:562:U:H1'	11:AL:11:ARG:HB3	1.65	0.78
1:AA:79:G:H2'	1:AA:80:A:C8	2.19	0.78
10:AK:92:ARG:HB3	10:AK:92:ARG:HH11	1.48	0.78
48:B1:12:SER:HA	48:B1:50:GLU:HB3	1.64	0.78
23:BB:597:G:H4'	33:BL:21:ARG:HE	1.48	0.78
34:BM:67:VAL:HB	34:BM:100:LYS:HG2	1.63	0.78
36:BO:37:ALA:O	36:BO:38:GLN:HG2	1.84	0.78
23:BB:328:U:H4'	42:BU:65:GLN:HG3	1.65	0.78
13:CN:40:ARG:HH12	18:CS:5:LYS:HB2	1.48	0.78
27:DE:137:LYS:HA	27:DE:137:LYS:NZ	1.98	0.78
52:DI:45:THR:HA	52:DI:48:ILE:HG22	1.64	0.78
35:DN:48:VAL:HA	35:DN:51:LEU:HD12	1.66	0.78
43:DW:66:VAL:HG13	43:DW:67:LYS:H	1.47	0.78
45:BY:16:LEU:HB2	45:BY:17:PRO:CD	2.14	0.78
23:DB:365:U:H2'	23:DB:366:C:C6	2.18	0.78
26:DD:170:VAL:HB	26:DD:194:PRO:HG2	1.66	0.78
26:DD:89:GLU:HG2	26:DD:93:GLY:HA3	1.64	0.78
37:DP:29:VAL:HG21	37:DP:61:ARG:HH22	1.48	0.78
3:AD:123:MET:HA	3:AD:128:VAL:HA	1.65	0.78
23:BB:146:A:H2'	23:BB:147:C:C6	2.19	0.78
34:BM:108:VAL:HG23	34:BM:112:LEU:HD11	1.65	0.78
3:AD:118:SER:HA	3:AD:130:ASN:HB2	1.65	0.78
3:AD:159:GLU:HG3	3:AD:160:LEU:N	1.96	0.78
27:BE:122:GLU:HB3	27:BE:152:GLU:HB2	1.65	0.78
5:CF:4:TYR:HA	5:CF:91:ARG:HA	1.64	0.78
8:CI:18:VAL:HG12	8:CI:64:ILE:HA	1.64	0.78
51:D4:26:ILE:O	51:D4:27:CYS:HB2	1.83	0.78
34:DM:134:THR:HG22	34:DM:135:VAL:H	1.46	0.78
34:DM:15:GLY:O	34:DM:16:ARG:HG3	1.83	0.78
13:AN:60:ARG:HE	13:AN:62:ARG:NE	1.82	0.78
23:BB:1283:G:N2	23:BB:1286:A:H5'	1.98	0.78
23:BB:483:A:H5'	42:BU:43:LYS:HE2	1.66	0.78
29:BG:173:ALA:C	29:BG:174:LYS:HD2	2.03	0.78
30:BH:37:VAL:N	30:BH:38:PRO:HD2	1.95	0.78
52:BI:25:PRO:O	52:BI:29:GLN:HG2	1.84	0.78
34:BM:40:ARG:HH11	34:BM:92:TRP:HE3	1.29	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BW:32:ALA:O	43:BW:66:VAL:HG11	1.84	0.78
10:CK:82:GLU:HB3	10:CK:108:ASN:HD22	1.48	0.78
21:CU:41:THR:HA	21:CU:45:LYS:HD3	1.65	0.78
37:DP:25:VAL:HG12	37:DP:27:VAL:H	1.47	0.78
24:DV:9:ARG:HG2	24:DV:41:GLU:HG2	1.66	0.78
45:DY:18:LYS:HD2	45:DY:18:LYS:N	1.99	0.78
23:BB:414:C:H2'	23:BB:415:A:H8	1.49	0.78
25:BC:21:PRO:HG2	25:BC:202:ARG:HH11	1.48	0.78
23:BB:2674:G:H4'	32:BK:30:ARG:HG2	1.66	0.78
42:BU:39:ASN:ND2	42:BU:42:LYS:HE2	1.99	0.78
9:CJ:40:ILE:HD11	9:CJ:73:LEU:HD12	1.65	0.78
23:DB:947:A:HO2'	23:DB:984:A:H2	1.30	0.78
23:DB:2511:U:H5''	26:DD:129:THR:CG2	2.14	0.78
23:DB:2811:G:H5'	26:DD:62:LYS:HZ1	1.46	0.78
28:DF:134:GLN:HB3	28:DF:136:ILE:HG13	1.65	0.78
31:DJ:41:LYS:NZ	38:DQ:63:ARG:HD2	1.99	0.78
35:DN:86:ARG:HB3	35:DN:86:ARG:HH11	1.47	0.78
20:AB:36:LYS:N	20:AB:36:LYS:HE3	1.98	0.78
5:AF:53:LYS:HA	5:AF:53:LYS:HZ3	1.49	0.78
12:AM:19:THR:HA	12:AM:24:VAL:HG23	1.65	0.78
23:BB:2867:G:H2'	23:BB:2867:G:N3	1.98	0.78
31:BJ:81:ILE:HG13	31:BJ:82:GLY:N	1.99	0.78
1:CA:17:U:H2'	1:CA:18:C:C6	2.19	0.78
1:CA:436:C:O2'	1:CA:437:U:H5'	1.83	0.78
5:CF:62:MET:HG3	5:CF:64:VAL:HG23	1.66	0.78
50:D3:7:ARG:HA	50:D3:7:ARG:HH11	1.49	0.78
45:DY:4:ILE:N	45:DY:36:GLU:HG3	1.99	0.78
18:AS:39:ILE:HG21	18:AS:65:MET:HB3	1.63	0.77
25:BC:127:ASN:ND2	25:BC:128:THR:H	1.82	0.77
38:BQ:33:VAL:HG23	38:BQ:37:ALA:HB3	1.66	0.77
25:DC:139:THR:HA	25:DC:193:GLU:CD	2.05	0.77
34:DM:117:PHE:HB2	34:DM:124:LEU:HD11	1.67	0.77
20:AB:58:LYS:HG2	20:AB:62:ARG:HH21	1.50	0.77
12:AM:23:GLY:HA3	12:AM:64:VAL:HG13	1.64	0.77
16:AQ:80:LYS:O	16:AQ:80:LYS:HE3	1.84	0.77
23:BB:663:G:H5''	33:BL:25:SER:HB2	1.64	0.77
1:CA:975:A:O2'	1:CA:1358:U:H1'	1.84	0.77
26:DD:77:ARG:HB2	26:DD:77:ARG:NH1	1.98	0.77
23:BB:918:A:H2'	23:BB:919:U:H5'	1.66	0.77
35:BN:116:VAL:HG22	35:BN:117:ASP:H	1.49	0.77
36:BO:38:GLN:HA	36:BO:50:ALA:HB3	1.67	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:25:VAL:HG12	37:BP:27:VAL:HG12	1.67	0.77
16:CQ:13:SER:HB3	16:CQ:21:VAL:HB	1.66	0.77
25:DC:229:HIS:ND1	25:DC:230:PRO:HD2	1.99	0.77
26:DD:8:LYS:HE2	37:DP:5:LYS:HG3	1.67	0.77
29:DG:151:ARG:HH11	29:DG:151:ARG:HB2	1.47	0.77
37:DP:90:ALA:H	37:DP:112:ARG:HH21	1.32	0.77
26:BD:106:LYS:HA	26:BD:176:ASP:HA	1.66	0.77
43:BW:43:LYS:HG2	43:BW:76:ARG:HH21	1.49	0.77
10:CK:23:HIS:HB3	10:CK:30:ILE:HG13	1.65	0.77
18:CS:27:LYS:HG3	18:CS:28:LYS:HD2	1.65	0.77
27:DE:50:ALA:O	27:DE:74:LYS:HD3	1.83	0.77
28:DF:174:PHE:H	28:DF:175:PRO:HD2	1.49	0.77
32:DK:76:VAL:HB	37:DP:74:GLN:HE21	1.50	0.77
1:AA:430:A:OP1	3:AD:8:LEU:HB2	1.84	0.77
23:BB:2578:G:H1'	26:BD:145:SER:OG	1.83	0.77
31:BJ:58:ASN:ND2	31:BJ:128:ASN:HB2	1.99	0.77
31:BJ:61:LYS:HE2	31:BJ:61:LYS:HA	1.64	0.77
34:BM:20:LEU:HD13	34:BM:38:ARG:HG2	1.66	0.77
13:CN:26:LEU:HA	13:CN:29:ILE:HD12	1.65	0.77
19:CT:69:ASN:H	19:CT:69:ASN:ND2	1.82	0.77
23:DB:1283:G:N2	23:DB:1286:A:H5'	1.98	0.77
51:B4:23:ILE:HG13	51:B4:35:GLN:HB3	1.65	0.77
23:BB:2576:G:OP2	26:BD:149:ASN:HB2	1.84	0.77
28:BF:126:ASN:HD22	28:BF:126:ASN:H	1.30	0.77
38:BQ:94:LEU:H	38:BQ:94:LEU:HD13	1.49	0.77
23:DB:966:G:H5'	23:DB:2272:U:O2	1.84	0.77
23:DB:2867:G:N3	23:DB:2867:G:H2'	1.98	0.77
8:AI:11:ARG:HH21	8:AI:12:LYS:HB2	1.49	0.77
30:BH:81:ALA:HA	30:BH:146:VAL:HA	1.66	0.77
34:BM:100:LYS:HE3	34:BM:100:LYS:H	1.49	0.77
52:DI:27:LEU:HD12	52:DI:32:VAL:HG11	1.66	0.77
39:DR:60:LYS:O	39:DR:60:LYS:HD3	1.85	0.77
41:DT:38:ALA:HB1	41:DT:43:ILE:HD11	1.65	0.77
1:AA:1412:C:H2'	1:AA:1413:A:C8	2.20	0.77
2:AC:58:ARG:HG2	2:AC:63:ILE:HG22	1.67	0.77
23:BB:2008:C:H2'	23:BB:2009:A:H8	1.50	0.77
26:BD:13:ARG:HA	26:BD:13:ARG:NH1	2.00	0.77
26:BD:62:LYS:HG2	26:BD:63:PRO:HD3	1.65	0.77
35:BN:100:CYS:HB3	35:BN:110:MET:HB2	1.66	0.77
23:DB:2293:G:H5''	36:DO:9:ARG:NH2	1.99	0.77
45:DY:45:GLY:HA2	45:DY:48:ASN:ND2	2.00	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AF:10:VAL:HA	5:AF:84:VAL:HA	1.67	0.77
49:B2:12:ARG:HH21	49:B2:18:PHE:H	1.31	0.77
23:BB:2064:C:H2'	23:BB:2065:C:C6	2.19	0.77
30:BH:73:ASN:ND2	30:BH:74:ALA:H	1.82	0.77
23:BB:2324:U:H5'	23:BB:2325:G:H5''	1.67	0.77
23:BB:441:U:O2'	23:BB:442:G:H5'	1.84	0.77
25:BC:153:LEU:HD23	25:BC:153:LEU:H	1.49	0.77
52:BI:21:PRO:HB2	52:BI:22:PRO:HD3	1.66	0.77
31:BJ:57:LEU:HB3	31:BJ:128:ASN:HA	1.66	0.77
23:DB:51:G:O2'	23:DB:118:A:N6	2.17	0.77
1:AA:1461:G:H2'	1:AA:1462:C:C6	2.20	0.76
50:B3:48:MET:SD	50:B3:51:LYS:HD3	2.25	0.76
23:DB:458:G:C5'	49:D2:39:ARG:HB2	2.13	0.76
23:DB:1486:U:H2'	23:DB:1487:U:C6	2.21	0.76
23:DB:2071:A:H2'	23:DB:2072:C:C6	2.20	0.76
23:DB:526:A:N6	23:DB:2626:C:H4'	2.00	0.76
52:DI:25:PRO:O	52:DI:29:GLN:HG3	1.84	0.76
15:AP:51:ARG:HB3	15:AP:51:ARG:HH11	1.48	0.76
31:BJ:7:LYS:HD3	31:BJ:48:VAL:HB	1.67	0.76
39:BR:47:VAL:HG22	39:BR:48:LYS:H	1.48	0.76
11:CL:78:VAL:HG12	11:CL:101:LEU:HD13	1.67	0.76
23:DB:2886:A:C8	47:D0:27:LEU:HG	2.21	0.76
23:DB:1826:G:OP2	25:DC:221:GLY:HA2	1.86	0.76
26:DD:8:LYS:HD3	37:DP:5:LYS:HE3	1.67	0.76
1:AA:1123:U:O2'	1:AA:1124:G:H5'	1.85	0.76
20:AB:202:ASN:ND2	20:AB:204:ASP:H	1.83	0.76
1:AA:1057:G:H5''	2:AC:153:SER:HB3	1.67	0.76
25:BC:106:PRO:HD2	25:BC:109:LEU:HD22	1.68	0.76
30:BH:3:VAL:HG22	30:BH:21:VAL:HG11	1.67	0.76
4:CE:110:MET:SD	4:CE:126:ALA:HB2	2.25	0.76
8:CI:118:ARG:HH21	8:CI:122:ARG:NE	1.82	0.76
9:CJ:57:VAL:HG22	9:CJ:58:ASN:H	1.49	0.76
23:DB:1113:U:H5''	29:DG:2:ARG:CZ	2.16	0.76
26:DD:5:VAL:HB	26:DD:27:ILE:O	1.85	0.76
37:DP:76:HIS:H	37:DP:76:HIS:HD2	1.30	0.76
24:DV:30:ILE:HG12	24:DV:91:PHE:HB2	1.65	0.76
10:AK:111:ASP:HB2	21:AU:16:ARG:HH22	1.50	0.76
21:AU:17:ARG:HA	21:AU:20:ARG:HB3	1.66	0.76
25:BC:224:MET:HA	25:BC:232:GLY:HA2	1.67	0.76
36:BO:19:GLN:H	43:BW:76:ARG:HH12	1.32	0.76
3:CD:138:PRO:HA	3:CD:181:PHE:HD2	1.50	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2008:C:H2'	23:DB:2009:A:H8	1.51	0.76
27:DE:14:VAL:HG12	27:DE:15:SER:H	1.48	0.76
7:AH:30:LYS:HA	7:AH:30:LYS:HZ2	1.49	0.76
11:AL:105:GLY:HA3	11:AL:117:GLY:HA3	1.67	0.76
11:AL:78:VAL:HG12	11:AL:101:LEU:HD23	1.67	0.76
23:BB:811:U:C5	33:BL:29:LYS:HD3	2.20	0.76
4:CE:98:ALA:HB2	4:CE:123:LEU:HG	1.68	0.76
1:CA:1225:A:H5'	12:CM:101:THR:OG1	1.84	0.76
38:DQ:30:VAL:HG12	38:DQ:31:TYR:N	2.00	0.76
1:AA:1251:A:H2'	1:AA:1252:A:C8	2.21	0.76
47:B0:6:LYS:HA	47:B0:6:LYS:NZ	1.99	0.76
23:BB:1022:G:N7	31:BJ:68:LYS:HE3	2.01	0.76
23:BB:1252:G:H1	38:BQ:36:GLN:NE2	1.83	0.76
23:BB:27:G:H22	23:BB:512:G:H2'	1.50	0.76
25:BC:77:VAL:HG23	25:BC:115:ILE:HD11	1.66	0.76
1:CA:939:G:H5''	6:CG:101:ARG:HH12	1.49	0.76
47:D0:41:HIS:CG	47:D0:42:ILE:H	2.03	0.76
23:DB:1368:G:C5'	49:D2:25:LYS:HG2	2.15	0.76
34:DM:20:LEU:HD13	34:DM:38:ARG:HG3	1.67	0.76
4:AE:131:ASN:HD21	4:AE:133:ILE:HB	1.50	0.76
48:B1:42:VAL:HG12	48:B1:43:ARG:HG3	1.67	0.76
23:BB:2420:C:OP1	50:B3:34:LYS:HB2	1.86	0.76
23:BB:2443:C:H2'	23:BB:2444:G:H8	1.51	0.76
25:BC:9:SER:HB2	25:BC:202:ARG:NE	2.00	0.76
27:BE:4:VAL:HG22	27:BE:5:LEU:HD12	1.65	0.76
40:BS:73:LYS:HB2	40:BS:106:VAL:HB	1.67	0.76
1:CA:814:A:H5'	1:CA:1511:G:H4'	1.66	0.76
4:CE:47:PHE:HE1	4:CE:137:ARG:HE	1.34	0.76
28:DF:132:ARG:HD3	28:DF:133:GLU:H	1.48	0.76
31:DJ:135:GLN:NE2	31:DJ:138:GLN:H	1.83	0.76
46:DZ:41:HIS:CG	46:DZ:42:PRO:HD2	2.21	0.76
23:BB:643:A:C2	48:B1:42:VAL:HG13	2.21	0.76
23:BB:1241:A:H2'	23:BB:1242:U:H5'	1.67	0.76
23:BB:125:A:H3'	23:BB:126:A:C5'	2.15	0.76
35:BN:119:SER:HB2	35:BN:121:LYS:HG2	1.67	0.76
37:BP:51:ASN:HA	37:BP:61:ARG:H	1.51	0.76
11:CL:49:ARG:HG2	11:CL:89:LEU:HD21	1.66	0.76
23:DB:2346:A:H3'	23:DB:2347:C:H5''	1.68	0.76
23:DB:547:A:H2'	23:DB:547:A:N3	1.99	0.76
26:DD:116:LYS:HG2	26:DD:123:LYS:HE2	1.67	0.76
30:DH:94:ILE:HG22	30:DH:122:LEU:HG	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DQ:48:ASP:HA	38:DQ:51:GLN:NE2	2.00	0.76
24:DV:9:ARG:HH12	24:DV:12:GLN:HA	1.49	0.76
8:AI:71:ILE:HD12	8:AI:71:ILE:H	1.50	0.76
28:BF:56:LEU:O	28:BF:59:ILE:HG22	1.86	0.76
30:BH:73:ASN:N	30:BH:73:ASN:HD22	1.83	0.76
52:BI:105:LEU:HD11	52:BI:139:VAL:HG11	1.66	0.76
9:CJ:25:ILE:HG23	9:CJ:87:LEU:HD12	1.68	0.76
23:DB:1098:A:P	52:DI:3:LYS:HG2	2.25	0.76
23:DB:1175:A:H2'	23:DB:1176:U:H5'	1.68	0.76
33:DL:19:LEU:H	33:DL:19:LEU:HD22	1.51	0.76
20:AB:41:ASN:ND2	20:AB:44:LYS:HB2	2.00	0.76
8:AI:33:SER:HB3	8:AI:36:GLN:HB2	1.68	0.76
10:AK:88:PRO:HD3	21:AU:28:LEU:HD13	1.68	0.76
12:AM:44:ILE:HD12	12:AM:45:SER:H	1.51	0.76
16:AQ:16:MET:HB2	16:AQ:19:SER:HB2	1.66	0.76
23:BB:118:A:H5'	23:BB:119:A:H8	1.49	0.76
25:BC:127:ASN:HD22	25:BC:128:THR:H	1.31	0.76
36:BO:76:LYS:HG2	36:BO:110:ALA:HB2	1.68	0.76
44:BX:45:GLN:H	44:BX:45:GLN:NE2	1.83	0.76
1:CA:1071:C:H2'	1:CA:1072:G:H8	1.52	0.76
7:CH:72:GLU:H	7:CH:129:ALA:HB2	1.51	0.76
23:DB:974:G:H1'	23:DB:975:A:C8	2.21	0.76
26:DD:60:VAL:HG23	26:DD:63:PRO:HG2	1.66	0.76
2:AC:8:GLY:HA2	2:AC:11:LEU:HG	1.69	0.75
2:AC:156:LEU:HD11	2:AC:163:ARG:O	1.87	0.75
2:AC:146:LYS:HE3	2:AC:202:PHE:HE2	1.51	0.75
3:AD:137:SER:HB2	3:AD:138:PRO:HD2	1.68	0.75
16:AQ:45:VAL:HG12	16:AQ:46:HIS:H	1.49	0.75
50:B3:12:ARG:HH11	50:B3:12:ARG:HB2	1.48	0.75
23:BB:2301:C:H2'	23:BB:2302:U:C6	2.21	0.75
25:BC:229:HIS:HE2	25:BC:244:VAL:HG22	1.48	0.75
31:BJ:7:LYS:HD2	31:BJ:49:ASP:HB2	1.66	0.75
33:BL:91:ASP:O	33:BL:92:LEU:HB2	1.86	0.75
34:BM:42:THR:HG22	34:BM:45:GLN:HB2	1.67	0.75
35:BN:38:LEU:HB2	35:BN:109:PRO:HB2	1.68	0.75
42:BU:78:LYS:HE3	42:BU:96:LYS:HB2	1.66	0.75
49:D2:43:THR:O	49:D2:44:VAL:HG13	1.85	0.75
22:DA:47:C:OP1	36:DO:1:MET:HA	1.87	0.75
24:DV:63:ILE:N	24:DV:70:ILE:HD11	1.99	0.75
2:AC:49:ALA:HB1	2:AC:75:VAL:HG22	1.68	0.75
5:AF:92:THR:HG23	5:AF:93:LYS:H	1.51	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:633:A:O5'	23:BB:633:A:H8	1.70	0.75
23:BB:78:U:H2'	23:BB:79:C:C6	2.21	0.75
27:BE:141:MET:HG3	27:BE:143:LEU:H	1.52	0.75
33:BL:94:THR:HB	33:BL:103:ILE:HG12	1.68	0.75
20:CB:218:ALA:HA	20:CB:221:ARG:HG2	1.68	0.75
20:CB:94:ARG:N	20:CB:94:ARG:HE	1.84	0.75
49:D2:35:ARG:HH21	49:D2:42:LEU:HD12	1.51	0.75
23:DB:957:C:OP1	34:DM:76:LYS:HD3	1.86	0.75
30:DH:75:LEU:H	30:DH:75:LEU:HD23	1.51	0.75
1:AA:967:C:H3'	1:AA:968:A:C5'	2.15	0.75
3:AD:159:GLU:HG3	3:AD:160:LEU:H	1.50	0.75
23:BB:1654:A:OP1	35:BN:1:MET:HB2	1.86	0.75
23:BB:855:G:H21	43:BW:23:LYS:HD2	1.50	0.75
34:BM:71:LYS:HD2	34:BM:71:LYS:O	1.86	0.75
43:BW:23:LYS:HB2	43:BW:23:LYS:HZ2	1.51	0.75
23:DB:1299:G:H4'	23:DB:1301:A:H1'	1.69	0.75
23:DB:2012:G:OP1	40:DS:98:LYS:HD3	1.85	0.75
23:DB:2204:G:O5'	25:DC:149:LYS:HE3	1.87	0.75
23:DB:534:U:H5'	38:DQ:41:ALA:HA	1.66	0.75
26:DD:179:ARG:HB2	26:DD:188:LEU:HD22	1.68	0.75
31:DJ:64:VAL:HG21	31:DJ:90:GLU:OE1	1.86	0.75
35:DN:42:LYS:HE3	35:DN:45:ARG:HG3	1.67	0.75
41:DT:14:PRO:HA	41:DT:32:LEU:HA	1.67	0.75
41:DT:55:VAL:HG23	41:DT:87:LEU:N	2.01	0.75
47:B0:8:THR:HG23	47:B0:9:ARG:H	1.51	0.75
23:BB:2346:A:H3'	23:BB:2347:C:H5''	1.67	0.75
26:BD:8:LYS:HB3	37:BP:5:LYS:NZ	2.01	0.75
28:BF:102:LEU:O	28:BF:107:VAL:HG23	1.86	0.75
29:BG:51:PHE:HE2	29:BG:63:GLN:HE21	1.34	0.75
38:BQ:91:ARG:N	39:BR:10:LYS:HZ2	1.79	0.75
45:BY:10:ARG:NE	45:BY:31:ILE:HG13	2.01	0.75
1:CA:1056:U:H5'	2:CC:162:ALA:HB2	1.65	0.75
1:CA:1399:C:H4'	1:CA:1400:C:H5''	1.67	0.75
1:CA:1508:A:H2'	1:CA:1509:C:C6	2.22	0.75
1:CA:780:A:O2'	1:CA:781:A:H5''	1.87	0.75
7:CH:10:LEU:HD22	7:CH:74:ILE:HD11	1.68	0.75
23:DB:2443:C:H2'	23:DB:2444:G:H8	1.51	0.75
26:DD:37:VAL:HB	26:DD:46:ARG:HB2	1.68	0.75
42:DU:4:ILE:HG21	42:DU:27:VAL:HG13	1.68	0.75
1:AA:1409:C:H42	1:AA:1491:G:H1	1.33	0.75
1:AA:575:G:H4'	1:AA:576:C:O5'	1.84	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AH:30:LYS:HA	7:AH:30:LYS:NZ	2.01	0.75
48:B1:12:SER:HB2	48:B1:21:THR:O	1.85	0.75
23:BB:1197:G:H2'	23:BB:1198:U:H6	1.49	0.75
23:BB:1440:U:H2'	23:BB:1441:G:H8	1.51	0.75
23:BB:1486:U:H2'	23:BB:1487:U:C6	2.21	0.75
26:BD:1:MET:HB2	26:BD:84:LEU:HD22	1.68	0.75
27:BE:89:PRO:O	27:BE:90:GLN:HG2	1.85	0.75
52:BI:20:SER:HB3	52:BI:21:PRO:HD3	1.68	0.75
39:BR:58:VAL:HG22	39:BR:59:ILE:H	1.52	0.75
44:BX:22:LEU:HD22	44:BX:26:PHE:HB2	1.69	0.75
1:CA:817:C:H1'	1:CA:819:A:H5'	1.69	0.75
3:CD:2:ARG:NH1	3:CD:114:ARG:HD3	2.00	0.75
23:DB:898:C:O2'	23:DB:899:A:H5''	1.86	0.75
42:DU:3:LYS:HD3	42:DU:81:ARG:HH12	1.52	0.75
24:DV:72:VAL:HG12	24:DV:93:ARG:HA	1.68	0.75
1:AA:814:A:H5'	1:AA:1511:G:H4'	1.69	0.75
2:AC:153:SER:HB2	2:AC:196:GLY:H	1.51	0.75
19:AT:59:ARG:HH11	19:AT:59:ARG:HB2	1.51	0.75
23:BB:65:U:OP1	41:BT:76:ARG:HB2	1.87	0.75
26:BD:91:THR:HG23	26:BD:92:VAL:H	1.50	0.75
27:BE:120:VAL:H	27:BE:189:THR:HG22	1.50	0.75
28:BF:35:LEU:HA	28:BF:153:ILE:HA	1.67	0.75
1:CA:269:C:H2'	1:CA:270:A:H8	1.50	0.75
4:CE:44:ARG:HA	4:CE:71:ILE:O	1.85	0.75
23:DB:1825:U:H5'	25:DC:244:VAL:CG2	2.16	0.75
23:DB:544:C:H2'	23:DB:545:U:C4	2.21	0.75
30:DH:11:ASN:HD22	30:DH:20:ASN:HD22	1.33	0.75
1:AA:780:A:O2'	1:AA:781:A:H5''	1.87	0.75
7:AH:5:PRO:HB2	7:AH:32:LYS:HZ1	1.49	0.75
16:AQ:18:LYS:HG2	16:AQ:48:GLU:HA	1.68	0.75
31:BJ:45:THR:HB	38:BQ:63:ARG:NH2	2.01	0.75
1:CA:1412:C:H2'	1:CA:1413:A:C8	2.21	0.75
2:CC:182:ASP:HB2	2:CC:203:LYS:HE2	1.67	0.75
23:DB:2756:U:H1'	23:DB:2757:A:H5''	1.69	0.75
23:DB:718:A:H2'	23:DB:719:C:H5'	1.67	0.75
23:DB:90:U:H3'	23:DB:91:A:C5'	2.17	0.75
26:DD:24:VAL:HG11	26:DD:193:VAL:HG11	1.69	0.75
31:DJ:133:ALA:HA	31:DJ:136:GLN:HB2	1.66	0.75
37:DP:28:LYS:H	37:DP:86:LYS:HD3	1.52	0.75
1:AA:1111:A:HO2'	1:AA:1112:C:H6	1.30	0.75
1:AA:65:A:N3	1:AA:65:A:H2'	2.00	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:235:GLU:HG3	25:BC:237:ARG:HD3	1.67	0.75
42:BU:7:ASP:N	42:BU:24:VAL:HA	2.01	0.75
23:DB:704:G:O2'	23:DB:727:A:N6	2.19	0.75
23:DB:2204:G:H4'	25:DC:149:LYS:HG3	1.69	0.75
36:DO:72:ALA:O	36:DO:76:LYS:HE3	1.87	0.75
31:DJ:40:HIS:HA	38:DQ:69:ARG:HH12	1.50	0.75
1:AA:1326:U:H2'	1:AA:1327:C:C6	2.21	0.75
1:AA:269:C:H2'	1:AA:270:A:H8	1.50	0.75
6:AG:67:ASN:HD22	6:AG:127:ALA:HA	1.51	0.75
23:BB:742:A:H2'	23:BB:743:A:C8	2.21	0.75
52:BI:122:GLU:O	52:BI:126:ARG:HG3	1.87	0.75
34:BM:42:THR:HA	34:BM:92:TRP:HD1	1.51	0.75
25:DC:257:ARG:HA	25:DC:261:ARG:HD2	1.68	0.75
23:DB:1060:U:C5	52:DI:131:THR:HG22	2.22	0.75
32:DK:102:PRO:HB3	32:DK:121:GLU:HG2	1.69	0.75
33:DL:55:MET:HB3	33:DL:56:PRO:HD3	1.69	0.75
37:DP:47:ILE:HG22	37:DP:48:ALA:N	2.00	0.75
1:AA:1125:U:H2'	1:AA:1126:U:H5''	1.68	0.74
16:AQ:10:ARG:NH1	16:AQ:56:ASP:H	1.85	0.74
23:BB:508:A:H2'	23:BB:509:C:H5'	1.66	0.74
23:BB:590:A:H2'	23:BB:591:U:C6	2.22	0.74
25:BC:77:VAL:HG21	25:BC:109:LEU:HG	1.69	0.74
33:BL:78:ARG:HG2	33:BL:99:ASN:HB3	1.69	0.74
1:CA:1277:C:O2'	1:CA:1279:G:H1'	1.86	0.74
8:CI:27:ILE:HD13	8:CI:34:LEU:HD13	1.68	0.74
23:DB:1099:G:C8	52:DI:3:LYS:HB2	2.22	0.74
23:DB:590:A:H2'	23:DB:591:U:C6	2.22	0.74
27:DE:142:ALA:N	27:DE:185:LYS:HZ1	1.83	0.74
23:DB:2751:G:H5'	29:DG:3:VAL:CG2	2.16	0.74
35:DN:8:ARG:HD2	35:DN:46:ARG:HE	1.48	0.74
1:AA:1000:A:H2'	1:AA:1001:C:C6	2.22	0.74
20:AB:61:SER:HA	20:AB:224:ARG:HA	1.69	0.74
11:AL:35:ARG:HA	11:AL:35:ARG:HE	1.52	0.74
23:BB:1178:C:H2'	23:BB:1179:G:H8	1.51	0.74
23:BB:2471:A:O2'	23:BB:2472:G:H8	1.70	0.74
23:BB:394:C:H2'	23:BB:395:U:O4'	1.86	0.74
25:BC:110:LYS:HE3	25:BC:111:ALA:H	1.52	0.74
27:BE:52:VAL:HG22	27:BE:53:THR:H	1.52	0.74
23:BB:2840:C:H5''	35:BN:53:THR:HG21	1.69	0.74
41:BT:29:THR:HG22	41:BT:86:THR:HG22	1.69	0.74
1:CA:1316:G:H2'	1:CA:1318:A:OP2	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CB:18:GLN:O	20:CB:37:VAL:HG23	1.86	0.74
3:CD:57:LYS:HB2	3:CD:199:ILE:HB	1.68	0.74
15:CP:4:ILE:HG12	15:CP:21:VAL:HG22	1.67	0.74
49:D2:25:LYS:HD2	49:D2:25:LYS:N	2.01	0.74
23:DB:742:A:H2'	23:DB:743:A:C8	2.21	0.74
25:DC:15:VAL:HG13	25:DC:16:VAL:HG23	1.67	0.74
25:DC:193:GLU:O	25:DC:194:VAL:HG13	1.87	0.74
37:DP:32:VAL:HB	37:DP:80:VAL:HG13	1.69	0.74
23:BB:1009:A:O2'	38:BQ:61:ILE:HD11	1.86	0.74
23:BB:107:G:H21	23:BB:346:A:N6	1.84	0.74
23:BB:1727:C:H2'	23:BB:1728:C:C6	2.22	0.74
23:BB:2086:U:H2'	23:BB:2087:G:C8	2.22	0.74
23:BB:275:C:H2'	23:BB:276:U:O4'	1.86	0.74
26:BD:35:THR:HG22	26:BD:46:ARG:HH22	1.52	0.74
23:BB:1113:U:H5''	29:BG:1:SER:HA	1.69	0.74
46:BZ:53:THR:HA	46:BZ:56:ARG:HD2	1.68	0.74
20:CB:81:ASP:HA	20:CB:84:LEU:HD22	1.68	0.74
3:CD:117:VAL:HG12	3:CD:130:ASN:HA	1.70	0.74
1:CA:878:A:O4'	7:CH:3:GLN:HG3	1.88	0.74
47:D0:32:THR:HG21	47:D0:41:HIS:NE2	2.01	0.74
49:D2:35:ARG:NH2	49:D2:43:THR:H	1.86	0.74
39:DR:4:VAL:HG12	39:DR:43:ASN:HB3	1.69	0.74
1:AA:1446:A:H2'	1:AA:1447:A:H5''	1.69	0.74
21:AU:40:PRO:HG2	21:AU:41:THR:H	1.51	0.74
23:BB:643:A:H2	48:B1:42:VAL:HG13	1.52	0.74
23:BB:857:G:O2'	23:BB:858:G:H5'	1.86	0.74
33:BL:118:THR:CG2	33:BL:119:PRO:HD3	2.17	0.74
33:BL:38:GLN:CD	33:BL:39:LYS:H	1.89	0.74
1:CA:1436:U:H2'	1:CA:1437:A:H8	1.52	0.74
1:CA:270:A:H2'	1:CA:271:C:C6	2.23	0.74
2:CC:95:GLY:O	2:CC:96:VAL:HG13	1.86	0.74
10:CK:110:THR:HG22	21:CU:5:VAL:H	1.52	0.74
30:DH:6:LEU:HD12	30:DH:36:ALA:N	2.02	0.74
36:DO:100:HIS:O	36:DO:104:GLN:HB3	1.87	0.74
38:DQ:7:VAL:HG13	38:DQ:8:ILE:HD12	1.67	0.74
11:AL:20:VAL:HG13	11:AL:94:TYR:HH	1.50	0.74
11:AL:22:ALA:HB1	11:AL:29:LYS:HG3	1.70	0.74
12:AM:78:ARG:NH2	12:AM:79:LEU:HG	2.03	0.74
25:BC:146:LYS:HB3	25:BC:147:PRO:HD2	1.69	0.74
25:BC:222:THR:HA	25:BC:231:HIS:NE2	2.02	0.74
30:BH:26:ALA:HB1	30:BH:31:VAL:HG23	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BM:8:LYS:HZ2	34:BM:70:ASP:H	1.34	0.74
35:BN:62:ASN:ND2	35:BN:80:PHE:HB3	2.02	0.74
35:BN:82:GLU:O	35:BN:86:ARG:HG3	1.88	0.74
38:BQ:106:THR:O	38:BQ:109:VAL:HG22	1.86	0.74
38:BQ:47:ARG:HH22	38:BQ:50:ARG:HB3	1.51	0.74
41:BT:5:GLU:HG2	41:BT:6:ARG:N	2.02	0.74
1:CA:430:A:OP1	3:CD:8:LEU:HB2	1.86	0.74
1:CA:1219:A:H5''	13:CN:52:ARG:HH12	1.52	0.74
14:CO:21:THR:HA	14:CO:26:VAL:HG11	1.69	0.74
16:CQ:5:ARG:HB3	16:CQ:7:LEU:HD23	1.70	0.74
42:DU:46:LYS:HE2	42:DU:53:GLN:HA	1.69	0.74
1:AA:1008:U:H2'	1:AA:1009:U:H4'	1.68	0.74
2:AC:179:ALA:HB1	2:AC:202:PHE:HE1	1.50	0.74
4:AE:87:VAL:HG12	4:AE:92:ARG:HA	1.68	0.74
34:BM:69:PRO:O	34:BM:70:ASP:HB3	1.86	0.74
45:BY:54:VAL:HG22	45:BY:55:LYS:H	1.52	0.74
23:DB:2102:G:H2'	23:DB:2103:C:O4'	1.87	0.74
23:DB:2144:G:O2'	23:DB:2145:C:H5'	1.87	0.74
30:DH:121:VAL:HG23	30:DH:122:LEU:HD23	1.68	0.74
40:DS:6:LYS:HA	40:DS:104:THR:HA	1.68	0.74
43:DW:42:THR:HG23	43:DW:66:VAL:H	1.53	0.74
23:BB:1437:C:H2'	23:BB:1438:U:C6	2.22	0.74
23:BB:19:A:H2'	23:BB:20:C:C6	2.23	0.74
27:BE:47:LYS:CD	27:BE:47:LYS:H	1.98	0.74
33:BL:77:ILE:HG22	33:BL:78:ARG:HG3	1.69	0.74
44:BX:22:LEU:HD22	44:BX:26:PHE:H	1.52	0.74
1:CA:1170:A:H2'	1:CA:1171:A:O4'	1.88	0.74
49:D2:25:LYS:HD2	49:D2:25:LYS:H	1.52	0.74
43:DW:3:LYS:HA	43:DW:3:LYS:HZ3	1.50	0.74
20:AB:120:SER:HA	20:AB:125:PHE:HB3	1.68	0.74
13:AN:42:ASN:HB3	13:AN:46:LYS:HE2	1.70	0.74
22:BA:60:C:H2'	22:BA:61:G:H8	1.53	0.74
23:BB:2420:C:OP1	50:B3:34:LYS:HE3	1.87	0.74
32:BK:11:ALA:HB3	32:BK:85:VAL:CG2	2.17	0.74
33:BL:110:VAL:HB	33:BL:126:ARG:NH2	2.02	0.74
34:BM:95:LEU:HD22	34:BM:98:PRO:HG3	1.69	0.74
42:BU:81:ARG:HH11	42:BU:81:ARG:HA	1.52	0.74
43:BW:48:ALA:HB3	43:BW:72:GLY:HA3	1.68	0.74
18:CS:50:VAL:HG21	18:CS:70:LEU:HG	1.70	0.74
50:D3:51:LYS:HA	50:D3:51:LYS:NZ	2.03	0.74
23:DB:1063:G:H1'	52:DI:92:PRO:HG2	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:704:G:H1'	23:DB:727:A:N6	2.03	0.74
27:DE:99:LYS:HZ2	27:DE:99:LYS:C	1.91	0.74
23:DB:2257:U:H5'	43:DW:5:ALA:HB2	1.69	0.74
1:AA:1294:G:H2'	1:AA:1295:U:C6	2.22	0.74
9:AJ:92:LEU:H	9:AJ:92:LEU:HD22	1.52	0.74
48:B1:53:ILE:HD13	48:B1:54:LYS:N	2.03	0.74
49:B2:13:ASN:HA	49:B2:18:PHE:HB2	1.69	0.74
23:BB:192:C:H2'	23:BB:193:U:H5'	1.69	0.74
26:BD:110:THR:HG22	26:BD:171:THR:HA	1.70	0.74
26:BD:13:ARG:HA	26:BD:13:ARG:HH11	1.53	0.74
26:BD:107:VAL:HA	26:BD:205:PRO:O	1.87	0.74
23:BB:37:C:C2'	27:BE:46:GLN:HG2	2.17	0.74
4:CE:82:HIS:HB2	4:CE:83:PRO:HD2	1.70	0.74
8:CI:4:GLN:HG2	8:CI:21:LYS:HD3	1.69	0.74
22:DA:17:C:H2'	22:DA:18:G:O4'	1.88	0.74
23:DB:1081:U:H5'	52:DI:126:ARG:NH1	2.02	0.74
23:DB:1785:A:H2'	23:DB:1787:A:N7	2.03	0.74
23:DB:633:A:O5'	23:DB:633:A:H8	1.69	0.74
29:DG:155:PRO:HB3	29:DG:171:LYS:HB3	1.68	0.74
52:DI:20:SER:HB3	52:DI:21:PRO:HD3	1.68	0.74
36:DO:21:LEU:HD23	36:DO:22:GLY:H	1.51	0.74
1:AA:270:A:H2'	1:AA:271:C:C6	2.23	0.74
1:AA:817:C:H1'	1:AA:819:A:H5'	1.69	0.74
13:AN:30:ILE:HG21	13:AN:41:TRP:HB2	1.68	0.74
15:AP:42:ILE:HG22	15:AP:43:ALA:H	1.52	0.74
23:BB:1176:U:H2'	23:BB:1177:G:O4'	1.88	0.74
23:BB:1252:G:O6	38:BQ:35:PHE:HB2	1.88	0.74
23:BB:467:G:OP1	49:B2:33:ARG:HD3	1.88	0.74
52:BI:7:TYR:HB2	52:BI:58:ILE:O	1.88	0.74
23:BB:588:U:OP1	33:BL:27:LEU:HB2	1.86	0.74
23:DB:1727:C:H2'	23:DB:1728:C:C6	2.22	0.74
23:DB:704:G:H1'	23:DB:727:A:H61	1.52	0.74
26:DD:33:ARG:HH11	26:DD:33:ARG:HB2	1.51	0.74
1:AA:501:C:H2'	1:AA:502:A:H8	1.52	0.73
23:BB:2301:C:H2'	23:BB:2302:U:H6	1.52	0.73
37:BP:62:LYS:HE3	37:BP:74:GLN:HE22	1.53	0.73
1:CA:239:U:OP1	1:CA:239:U:H4'	1.88	0.73
1:CA:501:C:H2'	1:CA:502:A:H8	1.52	0.73
1:CA:812:G:H2'	1:CA:812:G:N3	2.02	0.73
23:DB:143:C:O2	41:DT:3:ARG:HD2	1.88	0.73
23:DB:643:A:N6	23:DB:2370:G:H1'	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:6:LYS:HG2	27:DE:119:ILE:HB	1.69	0.73
24:DV:63:ILE:H	24:DV:70:ILE:CD1	1.99	0.73
2:AC:185:THR:HG22	2:AC:186:SER:H	1.52	0.73
51:B4:25:VAL:HG23	51:B4:35:GLN:HB2	1.68	0.73
23:BB:2239:G:H5'	25:BC:246:PRO:CD	2.19	0.73
23:BB:2511:U:H4'	26:BD:128:ARG:NH2	2.03	0.73
25:BC:107:LYS:CB	25:BC:194:VAL:HG11	2.17	0.73
30:BH:84:ALA:HA	30:BH:90:LEU:HG	1.70	0.73
33:BL:110:VAL:H	33:BL:126:ARG:HB2	1.52	0.73
34:BM:114:ARG:HH21	34:BM:125:PRO:HA	1.53	0.73
40:BS:34:ASP:N	40:BS:37:THR:HG22	2.03	0.73
1:CA:619:U:H3	3:CD:130:ASN:ND2	1.87	0.73
6:CG:58:LEU:H	6:CG:58:LEU:HD23	1.51	0.73
8:CI:114:LYS:HB2	8:CI:117:LEU:HD12	1.70	0.73
23:DB:1437:C:H2'	23:DB:1438:U:C6	2.22	0.73
26:DD:189:VAL:HG12	26:DD:190:LYS:H	1.52	0.73
52:DI:121:ILE:H	52:DI:121:ILE:HD13	1.52	0.73
48:B1:45:HIS:O	48:B1:46:VAL:HG12	1.87	0.73
23:BB:1299:G:H4'	23:BB:1301:A:H1'	1.68	0.73
23:BB:1785:A:H2'	23:BB:1787:A:N7	2.03	0.73
26:BD:138:LEU:HD23	26:BD:138:LEU:H	1.52	0.73
23:BB:2575:C:H4'	26:BD:148:GLN:CA	2.17	0.73
29:BG:157:LYS:HG3	29:BG:159:LYS:HZ2	1.53	0.73
34:BM:82:MET:HB2	34:BM:84:LYS:HZ2	1.53	0.73
23:BB:1392:A:N6	41:BT:18:GLU:HG2	2.04	0.73
45:BY:15:ARG:N	45:BY:15:ARG:HE	1.87	0.73
45:BY:20:LYS:HA	45:BY:24:LEU:HB2	1.69	0.73
23:DB:1098:A:H2'	52:DI:3:LYS:C	2.08	0.73
23:DB:578:G:N2	38:DQ:32:ARG:HH21	1.86	0.73
26:DD:27:ILE:HD13	26:DD:28:GLU:N	2.03	0.73
34:DM:2:LEU:HD12	34:DM:2:LEU:H	1.53	0.73
35:DN:86:ARG:HB3	35:DN:86:ARG:NH1	2.02	0.73
2:AC:155:ARG:H	2:AC:162:ALA:HA	1.52	0.73
10:AK:113:THR:HG21	21:AU:28:LEU:HD11	1.71	0.73
22:BA:50:A:H5''	36:BO:68:LYS:HD2	1.71	0.73
25:BC:129:LEU:HD23	25:BC:130:PRO:HD2	1.70	0.73
33:BL:95:LEU:H	33:BL:95:LEU:HD12	1.53	0.73
38:BQ:65:ASN:HB2	38:BQ:75:TYR:HB2	1.69	0.73
41:BT:30:ILE:HG23	41:BT:31:VAL:N	2.03	0.73
43:BW:23:LYS:HB2	43:BW:23:LYS:NZ	2.04	0.73
1:CA:1241:G:H2'	1:CA:1242:G:H8	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:268:U:H2'	1:CA:269:C:C6	2.23	0.73
29:DG:5:LYS:HD3	29:DG:68:ARG:HB3	1.68	0.73
43:DW:35:ILE:HG21	43:DW:70:VAL:HG11	1.71	0.73
23:BB:1171:G:H2'	23:BB:1172:C:O4'	1.88	0.73
23:BB:1714:U:H3'	23:BB:1715:G:C5'	2.18	0.73
29:BG:26:LYS:HA	29:BG:31:GLU:HG2	1.70	0.73
36:BO:50:ALA:HB2	36:BO:83:LEU:HD11	1.70	0.73
24:BV:70:ILE:HG12	24:BV:71:LYS:H	1.51	0.73
1:CA:922:G:H2'	1:CA:923:A:C8	2.23	0.73
23:DB:1820:U:H3	25:DC:197:ALA:HB1	1.53	0.73
23:DB:2185:U:H2'	23:DB:2186:G:O4'	1.88	0.73
23:DB:2895:G:H2'	23:DB:2896:C:C6	2.24	0.73
34:DM:41:LEU:HD23	34:DM:67:VAL:HG21	1.71	0.73
1:AA:268:U:H2'	1:AA:269:C:C6	2.23	0.73
23:BB:2682:A:C2	26:BD:22:ILE:HD11	2.22	0.73
28:BF:135:ILE:HD13	28:BF:138:PRO:HA	1.69	0.73
35:BN:99:LYS:NZ	35:BN:99:LYS:HB3	2.03	0.73
13:CN:9:GLU:HA	13:CN:12:ARG:HB3	1.71	0.73
16:CQ:59:GLU:O	16:CQ:75:VAL:HG22	1.88	0.73
23:DB:179:C:H5''	49:D2:28:ARG:NH1	2.04	0.73
23:DB:2157:G:N3	23:DB:2157:G:H2'	2.04	0.73
25:DC:68:ARG:NH2	25:DC:127:ASN:HA	2.02	0.73
34:DM:92:TRP:HD1	34:DM:93:VAL:H	1.36	0.73
36:DO:25:ARG:HE	36:DO:94:ARG:HH12	1.36	0.73
23:DB:2331:G:H4'	43:DW:69:GLU:HB2	1.71	0.73
1:AA:278:G:H21	1:AA:279:A:H62	1.36	0.73
23:DB:1406:U:H2'	23:DB:1407:G:H8	1.53	0.73
23:DB:2064:C:H2'	23:DB:2065:C:C6	2.24	0.73
14:CO:88:ARG:HH22	23:DB:714:U:H3'	1.54	0.73
23:DB:920:A:H2'	23:DB:921:C:H6	1.52	0.73
35:DN:34:ILE:HG22	35:DN:35:LYS:H	1.53	0.73
42:DU:59:GLU:HG3	42:DU:62:ALA:HB2	1.69	0.73
1:AA:1026:G:H2'	1:AA:1027:C:C6	2.24	0.73
1:AA:239:U:H4'	1:AA:239:U:OP1	1.88	0.73
11:AL:8:ARG:HG3	11:AL:9:LYS:N	2.00	0.73
26:BD:159:LYS:HB2	26:BD:159:LYS:NZ	2.04	0.73
34:BM:8:LYS:HE3	34:BM:70:ASP:HA	1.71	0.73
20:CB:59:ILE:HD12	20:CB:60:ALA:N	2.04	0.73
13:CN:30:ILE:HG22	13:CN:41:TRP:HB3	1.70	0.73
22:DA:32:U:H1'	22:DA:52:A:N7	2.03	0.73
23:DB:2684:U:H4'	32:DK:76:VAL:HG21	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:149:ILE:HG12	27:DE:186:VAL:HA	1.71	0.73
28:DF:56:LEU:HA	28:DF:59:ILE:HG22	1.71	0.73
34:DM:5:LYS:HB2	34:DM:69:PRO:HG2	1.71	0.73
41:DT:67:VAL:HG12	41:DT:68:LYS:H	1.53	0.73
42:DU:24:VAL:HG12	42:DU:26:ASN:OD1	1.87	0.73
1:AA:1342:C:H2'	1:AA:1343:G:H8	1.54	0.73
1:AA:812:G:H2'	1:AA:812:G:N3	2.02	0.73
12:AM:11:HIS:N	12:AM:44:ILE:HD11	2.03	0.73
18:AS:48:ILE:HB	18:AS:59:VAL:HB	1.71	0.73
51:B4:24:ARG:HH22	51:B4:36:ARG:HD2	1.54	0.73
30:BH:73:ASN:HD22	30:BH:73:ASN:H	1.33	0.73
1:CA:278:G:H21	1:CA:279:A:H62	1.36	0.73
1:CA:575:G:H4'	1:CA:576:C:O5'	1.87	0.73
20:CB:130:LYS:HB3	20:CB:134:LEU:HB2	1.70	0.73
16:CQ:11:VAL:HG13	16:CQ:20:ILE:HG23	1.71	0.73
23:DB:1440:U:H2'	23:DB:1441:G:H8	1.51	0.73
27:DE:49:ARG:HG3	27:DE:52:VAL:CG2	2.18	0.73
29:DG:175:LYS:HG2	29:DG:176:LYS:H	1.54	0.73
52:DI:72:THR:HG22	52:DI:115:ASP:OD2	1.89	0.73
52:DI:55:PRO:HD3	52:DI:74:PRO:HD3	1.69	0.73
38:DQ:43:GLN:HA	39:DR:79:ARG:HG2	1.71	0.73
1:AA:1498:U:H3'	53:AA:1601:KSG:N1	2.04	0.73
22:BA:20:G:H2'	22:BA:21:G:H8	1.54	0.73
23:BB:222:A:N6	23:BB:232:G:H1'	2.03	0.73
23:BB:704:G:O2'	23:BB:727:A:N6	2.22	0.73
23:BB:934:U:H2'	23:BB:935:C:C6	2.24	0.73
27:BE:195:GLN:H	27:BE:199:MET:HA	1.54	0.73
34:BM:69:PRO:HB3	34:BM:93:VAL:HG12	1.70	0.73
20:CB:99:MET:HA	20:CB:106:VAL:HG21	1.71	0.73
23:DB:2814:A:H2'	23:DB:2815:C:H6	1.54	0.73
23:DB:950:G:H2'	23:DB:951:C:C6	2.24	0.73
34:DM:71:LYS:HA	34:DM:71:LYS:HZ3	1.53	0.73
42:DU:38:ILE:HD13	42:DU:64:ILE:HG13	1.70	0.73
45:DY:26:LEU:HD12	45:DY:28:LEU:HD22	1.70	0.73
8:AI:117:LEU:HD22	8:AI:123:ARG:HG2	1.71	0.72
22:BA:60:C:H2'	22:BA:61:G:C8	2.24	0.72
23:BB:1041:G:H2'	23:BB:1042:G:H8	1.53	0.72
23:BB:942:G:H5''	33:BL:40:SER:HB3	1.70	0.72
25:BC:140:VAL:O	25:BC:141:HIS:HB2	1.89	0.72
33:BL:79:LEU:HA	33:BL:113:ALA:HB3	1.71	0.72
34:BM:4:PRO:HD3	34:BM:47:GLU:OE1	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:27:VAL:HG21	37:BP:86:LYS:HB2	1.71	0.72
1:CA:1179:A:H2'	1:CA:1180:A:O4'	1.87	0.72
1:CA:492:C:H2'	1:CA:493:A:H5''	1.71	0.72
11:CL:17:LYS:HB2	11:CL:17:LYS:NZ	2.04	0.72
23:DB:1241:A:H2'	23:DB:1242:U:H5'	1.69	0.72
23:DB:631:A:HO2'	33:DL:66:PHE:HD1	1.37	0.72
25:DC:141:HIS:HB3	25:DC:190:THR:HB	1.69	0.72
23:DB:1099:G:OP2	52:DI:3:LYS:HA	1.89	0.72
33:DL:77:ILE:N	33:DL:77:ILE:HD12	2.03	0.72
42:DU:4:ILE:HD13	42:DU:5:ARG:N	2.04	0.72
42:DU:33:VAL:HG23	42:DU:65:GLN:NE2	2.04	0.72
1:AA:922:G:H4'	4:AE:24:VAL:HA	1.70	0.72
1:AA:927:G:H4'	1:AA:1503:A:N7	2.04	0.72
2:AC:53:ARG:HG2	2:AC:54:ILE:H	1.54	0.72
12:AM:18:LEU:HB3	12:AM:29:SER:HB2	1.69	0.72
25:BC:79:ARG:NH1	25:BC:80:LEU:H	1.87	0.72
43:BW:66:VAL:HG13	43:BW:67:LYS:N	2.04	0.72
2:CC:61:LYS:HB3	2:CC:61:LYS:HZ3	1.54	0.72
4:CE:44:ARG:NH1	4:CE:72:ASN:HD21	1.86	0.72
22:DA:51:G:C2'	22:DA:52:A:H5''	2.19	0.72
23:DB:1714:U:H3'	23:DB:1715:G:C5'	2.18	0.72
30:DH:39:ALA:O	30:DH:43:ASN:HB2	1.89	0.72
38:DQ:69:ARG:HA	38:DQ:73:ILE:HG22	1.69	0.72
39:DR:47:VAL:HG12	39:DR:49:ILE:HG12	1.71	0.72
1:AA:78:A:H2'	1:AA:79:G:C8	2.24	0.72
9:AJ:56:HIS:O	9:AJ:57:VAL:HG12	1.89	0.72
23:BB:459:U:H5'	49:B2:41:ARG:HE	1.53	0.72
23:BB:2722:G:H4'	35:BN:4:ARG:HB3	1.69	0.72
23:BB:2895:G:H2'	23:BB:2896:C:C6	2.23	0.72
41:BT:12:ARG:HE	44:BX:29:ARG:HH22	1.35	0.72
1:CA:120:A:H2'	1:CA:121:U:H5''	1.71	0.72
1:CA:1111:A:N1	2:CC:176:THR:HG22	2.03	0.72
23:DB:151:C:H2'	23:DB:152:A:H8	1.52	0.72
23:DB:1813:G:H21	25:DC:50:THR:HG23	1.55	0.72
33:DL:92:LEU:HD22	33:DL:96:LYS:HG2	1.72	0.72
41:DT:62:VAL:HG23	41:DT:63:VAL:H	1.55	0.72
43:DW:69:GLU:HG3	43:DW:70:VAL:H	1.52	0.72
1:AA:1062:U:H2'	1:AA:1063:C:C6	2.24	0.72
1:AA:390:U:H2'	1:AA:391:G:C8	2.25	0.72
48:B1:14:ALA:HA	48:B1:48:TYR:HA	1.71	0.72
22:BA:48:U:H2'	22:BA:49:C:C6	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BG:9:VAL:HG23	29:BG:47:ASN:HB3	1.70	0.72
46:BZ:3:LYS:HE3	46:BZ:8:LYS:HA	1.69	0.72
1:CA:390:U:H2'	1:CA:391:G:C8	2.25	0.72
5:CF:11:HIS:ND1	5:CF:12:PRO:HD2	2.04	0.72
6:CG:90:VAL:HG13	6:CG:94:ARG:HD3	1.69	0.72
13:CN:82:LYS:HE2	13:CN:85:GLU:HG3	1.71	0.72
19:CT:38:ILE:HD11	19:CT:82:ILE:HG22	1.71	0.72
47:D0:36:LYS:HB2	47:D0:41:HIS:HA	1.72	0.72
23:DB:2471:A:O2'	23:DB:2472:G:H8	1.71	0.72
27:DE:153:LEU:HD13	27:DE:154:ASP:N	2.03	0.72
27:DE:21:ARG:HH12	27:DE:23:PHE:HB3	1.55	0.72
52:DI:42:ASN:HA	52:DI:45:THR:OG1	1.90	0.72
7:AH:76:ARG:HG2	7:AH:77:VAL:H	1.55	0.72
23:BB:1774:C:H2'	23:BB:1774:C:O2	1.88	0.72
23:BB:588:U:H5'	33:BL:28:GLY:HA2	1.71	0.72
23:BB:616:A:H3'	23:BB:617:G:H8	1.54	0.72
23:BB:2575:C:H5''	26:BD:149:ASN:CB	2.19	0.72
23:BB:811:U:H2'	33:BL:31:GLY:CA	2.18	0.72
6:CG:72:VAL:HG12	6:CG:89:GLU:HA	1.71	0.72
1:CA:1060:U:H4'	9:CJ:54:SER:HB2	1.70	0.72
13:CN:29:ILE:HG22	13:CN:30:ILE:HG13	1.71	0.72
23:DB:974:G:H1'	23:DB:975:A:H8	1.54	0.72
25:DC:143:VAL:HG12	25:DC:144:GLU:H	1.52	0.72
24:DV:7:GLU:HA	24:DV:65:VAL:HG23	1.71	0.72
1:AA:613:C:OP1	3:AD:80:ARG:HG3	1.90	0.72
2:AC:26:LYS:HD3	2:AC:27:GLU:N	2.05	0.72
23:BB:1324:G:H1'	23:BB:1616:A:N6	2.05	0.72
23:BB:1406:U:H2'	23:BB:1407:G:H8	1.52	0.72
23:BB:2064:C:H2'	23:BB:2065:C:H6	1.54	0.72
23:BB:370:G:O2'	23:BB:423:A:H3'	1.89	0.72
27:BE:98:LYS:HG3	27:BE:99:LYS:N	2.05	0.72
31:BJ:20:ALA:HB1	31:BJ:59:ALA:HA	1.70	0.72
33:BL:21:ARG:NH1	33:BL:21:ARG:HB3	2.04	0.72
24:BV:24:ASN:O	24:BV:44:HIS:HB2	1.89	0.72
1:CA:269:C:H2'	1:CA:270:A:C8	2.24	0.72
50:D3:12:ARG:HB3	50:D3:23:HIS:CB	2.20	0.72
23:DB:1275:A:H2	23:DB:1645:G:H21	1.35	0.72
27:DE:147:LEU:HB3	27:DE:167:VAL:HG13	1.71	0.72
38:DQ:111:LYS:HE2	39:DR:52:PRO:HG3	1.71	0.72
43:DW:58:LEU:HG	43:DW:81:ILE:HA	1.71	0.72
1:AA:269:C:H2'	1:AA:270:A:C8	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AE:23:THR:HG23	4:AE:28:ARG:HD3	1.70	0.72
23:BB:1224:U:H4'	39:BR:90:ARG:HB3	1.71	0.72
23:BB:1275:A:H2	23:BB:1645:G:H21	1.35	0.72
23:BB:2328:A:H2'	23:BB:2329:U:C6	2.24	0.72
23:BB:2814:A:H2'	23:BB:2815:C:H6	1.54	0.72
31:BJ:62:VAL:HG21	31:BJ:101:ILE:HD11	1.72	0.72
32:BK:2:ILE:HD11	32:BK:8:LEU:HD11	1.71	0.72
23:BB:142:A:N3	41:BT:2:ILE:HD12	2.04	0.72
23:DB:2020:A:H5'	47:D0:8:THR:CG2	2.19	0.72
23:DB:2091:C:H3'	23:DB:2092:U:C5'	2.19	0.72
23:DB:2291:U:H2'	23:DB:2292:U:C6	2.25	0.72
23:DB:414:C:H2'	23:DB:415:A:C8	2.24	0.72
26:DD:61:THR:HG23	26:DD:62:LYS:HD3	1.70	0.72
27:DE:130:LYS:NZ	27:DE:130:LYS:HB2	2.05	0.72
35:DN:102:PHE:HD1	40:DS:40:ASN:HD21	1.37	0.72
37:DP:49:ILE:O	37:DP:50:ARG:HD3	1.89	0.72
6:AG:148:LYS:HG3	6:AG:151:ALA:HB3	1.71	0.72
13:AN:60:ARG:HH21	13:AN:62:ARG:HE	1.36	0.72
9:AJ:67:ILE:HG13	13:AN:95:LEU:HD13	1.71	0.72
23:BB:155:A:H2'	23:BB:156:A:C8	2.24	0.72
23:BB:2241:A:H2'	23:BB:2242:G:C8	2.25	0.72
26:BD:106:LYS:NZ	26:BD:208:LYS:HD3	2.04	0.72
27:BE:147:LEU:O	27:BE:148:ILE:HG13	1.90	0.72
34:BM:3:GLN:HG2	34:BM:46:ILE:HB	1.70	0.72
37:BP:26:GLU:HA	37:BP:46:VAL:HG13	1.71	0.72
46:BZ:8:LYS:CD	46:BZ:9:TYR:H	2.01	0.72
1:CA:1000:A:H2'	1:CA:1001:C:C6	2.23	0.72
14:CO:38:LEU:HD23	14:CO:55:LEU:HD13	1.72	0.72
16:CQ:68:LYS:HG2	16:CQ:69:THR:HG23	1.72	0.72
23:DB:1774:C:H2'	23:DB:1774:C:O2	1.88	0.72
23:DB:2898:U:O2'	31:DJ:137:PRO:HB3	1.89	0.72
23:DB:836:G:H2'	23:DB:837:C:C6	2.24	0.72
25:DC:10:PRO:HD2	25:DC:202:ARG:HH22	1.53	0.72
23:DB:833:A:H1'	33:DL:54:GLN:OE1	1.89	0.72
33:DL:77:ILE:HD12	33:DL:77:ILE:H	1.54	0.72
35:DN:56:LYS:HG3	35:DN:57:THR:H	1.55	0.72
1:AA:1308:U:H2'	1:AA:1309:G:H8	1.55	0.72
1:AA:859:G:H2'	1:AA:860:A:C8	2.25	0.72
1:AA:95:C:H2'	1:AA:95:C:O2	1.87	0.72
13:AN:12:ARG:HA	13:AN:15:LEU:HD11	1.71	0.72
23:BB:1275:A:N3	23:BB:1275:A:H3'	2.05	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:136:G:H2'	23:BB:137:U:C6	2.25	0.72
35:BN:28:LEU:HD11	35:BN:44:LEU:HD21	1.72	0.72
1:CA:439:U:H4'	3:CD:120:LYS:HD2	1.72	0.72
23:DB:1309:G:H5''	49:D2:9:VAL:HG13	1.72	0.72
28:DF:45:ASP:O	28:DF:46:LYS:HG3	1.90	0.72
35:DN:28:LEU:HA	35:DN:34:ILE:HD11	1.72	0.72
1:AA:1035:A:H2'	1:AA:1036:A:H8	1.53	0.72
6:AG:94:ARG:HD3	6:AG:98:LEU:HD11	1.72	0.72
23:BB:324:A:H2'	23:BB:325:G:O4'	1.89	0.72
27:BE:44:ARG:H	27:BE:44:ARG:HD3	1.53	0.72
33:BL:30:THR:HA	33:BL:36:LYS:HE3	1.71	0.72
2:CC:128:MET:SD	2:CC:131:ARG:HB2	2.29	0.72
4:CE:104:ILE:HD12	4:CE:111:ARG:HD3	1.72	0.72
4:CE:23:THR:HA	4:CE:28:ARG:HA	1.72	0.72
6:CG:12:LEU:HD22	6:CG:13:PRO:HD2	1.72	0.72
23:DB:1469:A:H2'	23:DB:1470:A:C8	2.25	0.72
23:DB:2328:A:H2'	23:DB:2329:U:C6	2.24	0.72
33:DL:109:LYS:HE2	33:DL:126:ARG:CZ	2.20	0.72
36:DO:28:VAL:HG13	36:DO:106:LEU:HD13	1.72	0.72
36:DO:25:ARG:HG2	36:DO:94:ARG:HH22	1.54	0.72
38:DQ:69:ARG:HB3	38:DQ:69:ARG:NH1	2.05	0.72
40:DS:42:LYS:O	40:DS:45:VAL:HG13	1.90	0.72
43:DW:43:LYS:O	43:DW:44:PHE:HB2	1.87	0.72
1:AA:275:G:C5'	16:AQ:15:LYS:HG2	2.20	0.71
2:AC:13:ILE:O	2:AC:14:VAL:HG22	1.90	0.71
5:AF:54:LEU:HD13	5:AF:55:HIS:H	1.55	0.71
7:AH:49:LYS:HG3	7:AH:50:VAL:H	1.55	0.71
23:BB:2246:G:H2'	23:BB:2247:A:H8	1.55	0.71
23:BB:2266:A:C4'	23:BB:2267:A:N7	2.53	0.71
26:BD:175:LEU:HD11	26:BD:192:ALA:HB3	1.72	0.71
27:BE:159:LEU:HA	27:BE:169:VAL:HG21	1.71	0.71
30:BH:57:LYS:HG3	30:BH:61:VAL:HG21	1.70	0.71
34:BM:37:GLY:H	34:BM:97:GLN:HE21	1.35	0.71
42:BU:60:LYS:HG3	42:BU:61:GLU:H	1.54	0.71
1:CA:1289:A:H3'	1:CA:1290:G:H8	1.55	0.71
8:CI:79:ARG:O	8:CI:79:ARG:HD2	1.90	0.71
23:DB:1118:C:H2'	23:DB:1119:U:H6	1.54	0.71
23:DB:1275:A:N3	23:DB:1275:A:H3'	2.05	0.71
23:DB:1429:G:H2'	23:DB:1430:G:H8	1.55	0.71
23:DB:289:G:H2'	23:DB:290:U:C6	2.24	0.71
25:DC:28:PRO:HG2	25:DC:79:ARG:CZ	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:598:U:H5'	33:DL:21:ARG:HB2	1.70	0.71
39:DR:65:ALA:HB3	39:DR:100:GLY:H	1.55	0.71
1:AA:545:C:O2'	1:AA:546:A:H5'	1.90	0.71
20:AB:67:LEU:H	20:AB:67:LEU:HD22	1.54	0.71
4:AE:105:ILE:HB	4:AE:123:LEU:HA	1.70	0.71
1:AA:1060:U:H5''	9:AJ:53:ILE:HG22	1.71	0.71
23:BB:2147:A:C5'	23:BB:2148:G:H5'	2.19	0.71
23:BB:2331:G:H4'	43:BW:69:GLU:HG2	1.70	0.71
23:BB:2461:A:H2'	23:BB:2462:C:C6	2.24	0.71
23:BB:532:A:H4'	23:BB:533:G:C8	2.25	0.71
30:BH:8:LYS:HD3	30:BH:8:LYS:O	1.90	0.71
39:BR:6:GLN:HG2	39:BR:10:LYS:HE2	1.72	0.71
41:BT:54:GLU:H	41:BT:91:GLN:NE2	1.86	0.71
1:CA:1361:G:H2'	1:CA:1362:A:H5''	1.73	0.71
13:CN:73:LEU:HD12	13:CN:83:VAL:HG21	1.70	0.71
23:DB:2039:U:H2'	23:DB:2040:G:H8	1.56	0.71
1:AA:120:A:H2'	1:AA:121:U:H5''	1.70	0.71
48:B1:31:GLU:HG3	48:B1:32:LYS:H	1.54	0.71
51:B4:25:VAL:HB	51:B4:33:HIS:O	1.90	0.71
23:BB:543:G:C2'	23:BB:544:C:H5''	2.19	0.71
25:BC:20:ASN:HB2	25:BC:203:VAL:HG13	1.72	0.71
26:BD:52:THR:HG22	26:BD:75:ALA:HB1	1.73	0.71
29:BG:175:LYS:HD3	29:BG:176:LYS:HB2	1.72	0.71
52:BI:89:SER:HA	52:BI:97:VAL:HG21	1.72	0.71
34:BM:71:LYS:HD3	34:BM:92:TRP:C	2.09	0.71
11:CL:98:ARG:HB3	11:CL:116:TYR:HA	1.71	0.71
48:D1:47:ILE:HG22	48:D1:48:TYR:N	2.03	0.71
23:DB:1175:A:C2'	23:DB:1176:U:H5'	2.19	0.71
23:DB:643:A:H61	23:DB:2370:G:H1'	1.55	0.71
23:DB:2461:A:H2'	23:DB:2462:C:C6	2.25	0.71
25:DC:68:ARG:HH22	25:DC:127:ASN:HA	1.54	0.71
28:DF:7:TYR:O	28:DF:11:VAL:HB	1.90	0.71
39:DR:18:GLN:HA	39:DR:99:THR:HA	1.73	0.71
20:AB:41:ASN:HD22	20:AB:44:LYS:HB2	1.53	0.71
3:AD:154:VAL:HG13	3:AD:155:LYS:H	1.55	0.71
26:BD:128:ARG:HA	26:BD:128:ARG:HE	1.55	0.71
33:BL:78:ARG:H	33:BL:110:VAL:HG13	1.55	0.71
1:CA:559:A:H4'	1:CA:560:A:H3'	1.73	0.71
3:CD:63:ILE:HG23	3:CD:64:TYR:CD1	2.24	0.71
21:CU:15:LEU:HA	21:CU:17:ARG:HH11	1.56	0.71
22:DA:76:G:H2'	22:DA:77:U:H6	1.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:11:C:H2'	23:DB:12:U:H5'	1.72	0.71
23:DB:2784:U:H4'	26:DD:42:ASN:H	1.54	0.71
23:DB:438:G:H2'	23:DB:439:A:H8	1.56	0.71
23:DB:483:A:H2'	23:DB:484:C:H5'	1.70	0.71
26:DD:37:VAL:HG11	26:DD:46:ARG:HD3	1.72	0.71
23:DB:448:U:H3'	27:DE:79:ARG:HH21	1.55	0.71
52:DI:73:PRO:HG2	52:DI:78:LEU:HD21	1.72	0.71
23:DB:856:G:C4'	43:DW:23:LYS:HD2	2.19	0.71
1:AA:275:G:H5'	16:AQ:15:LYS:HG2	1.72	0.71
1:AA:1049:U:H2'	13:AN:2:LYS:HD3	1.70	0.71
50:B3:44:ARG:HG3	50:B3:45:PRO:HD2	1.73	0.71
23:BB:1395:A:H4'	23:BB:1397:U:C5	2.26	0.71
23:BB:1797:G:O3'	25:BC:253:GLY:HA2	1.91	0.71
28:BF:12:VAL:HG13	28:BF:27:VAL:HG11	1.71	0.71
33:BL:60:ARG:HA	33:BL:60:ARG:NH1	2.06	0.71
41:BT:88:LYS:HG2	41:BT:89:GLU:N	2.04	0.71
45:BY:9:THR:HG23	45:BY:10:ARG:H	1.55	0.71
1:CA:1009:U:H1'	1:CA:1021:A:N1	2.05	0.71
1:CA:84:U:O2'	1:CA:85:U:H5'	1.89	0.71
1:CA:8:A:H1'	4:CE:107:GLY:HA2	1.71	0.71
5:CF:38:ARG:HH21	5:CF:63:ASN:ND2	1.88	0.71
6:CG:14:ASP:HB3	6:CG:18:GLY:N	2.06	0.71
9:CJ:66:GLU:HB2	13:CN:98:ALA:HB2	1.72	0.71
17:CR:51:GLN:HA	17:CR:54:LEU:HD13	1.71	0.71
21:CU:14:ALA:H	21:CU:16:ARG:NH2	1.89	0.71
23:DB:1060:U:C4	23:DB:1088:A:N6	2.58	0.71
23:DB:1395:A:H4'	23:DB:1397:U:C5	2.26	0.71
23:DB:480:A:O4'	42:DU:40:LEU:HD22	1.89	0.71
25:DC:127:ASN:HD22	25:DC:128:THR:N	1.87	0.71
26:DD:6:GLY:HA2	26:DD:199:SER:O	1.90	0.71
33:DL:77:ILE:HD13	33:DL:110:VAL:N	2.05	0.71
38:DQ:46:TYR:HA	39:DR:77:PHE:HD2	1.56	0.71
45:DY:3:THR:HA	45:DY:37:ARG:H	1.56	0.71
1:AA:1208:C:H2'	1:AA:1209:C:O4'	1.91	0.71
1:AA:492:C:H2'	1:AA:493:A:H5''	1.71	0.71
1:AA:8:A:H2'	3:AD:205:LYS:O	1.91	0.71
49:B2:13:ASN:HB2	49:B2:18:PHE:HD2	1.56	0.71
25:BC:28:PRO:O	25:BC:29:PHE:HB3	1.90	0.71
29:BG:94:ARG:HB2	29:BG:94:ARG:HH11	1.55	0.71
39:BR:74:ILE:HA	39:BR:90:ARG:NE	2.03	0.71
23:DB:1324:G:H1'	23:DB:1616:A:N6	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:363:G:H2'	23:DB:364:C:C6	2.25	0.71
23:DB:616:A:H3'	23:DB:617:G:H8	1.54	0.71
26:DD:125:TRP:HD1	26:DD:127:PHE:HB2	1.56	0.71
1:AA:1284:C:H3'	1:AA:1285:A:H5''	1.72	0.71
1:AA:17:U:H2'	1:AA:18:C:H6	1.55	0.71
1:AA:188:C:H2'	1:AA:189:A:O4'	1.90	0.71
1:AA:559:A:H4'	1:AA:560:A:H3'	1.73	0.71
6:AG:38:ALA:O	6:AG:41:ILE:HG22	1.91	0.71
1:AA:1250:A:H4'	8:AI:69:GLY:H	1.56	0.71
10:AK:42:GLY:HA3	10:AK:73:VAL:HG22	1.71	0.71
12:AM:64:VAL:HG12	12:AM:65:GLU:H	1.54	0.71
51:B4:33:HIS:CD2	51:B4:33:HIS:H	2.07	0.71
51:B4:7:VAL:HG12	51:B4:8:LYS:H	1.54	0.71
22:BA:48:U:H2'	22:BA:49:C:H6	1.56	0.71
23:BB:1041:G:H2'	23:BB:1042:G:C8	2.25	0.71
23:BB:1173:U:C4	23:BB:1174:U:H1'	2.26	0.71
23:BB:1429:G:H2'	23:BB:1430:G:H8	1.55	0.71
26:BD:31:ALA:HB1	26:BD:49:GLN:HG3	1.73	0.71
28:BF:31:GLU:H	28:BF:157:THR:HA	1.54	0.71
33:BL:100:ILE:HG12	33:BL:102:GLY:H	1.55	0.71
35:BN:35:LYS:HA	35:BN:111:ALA:O	1.91	0.71
43:BW:19:ARG:NH2	43:BW:71:LYS:HD3	2.06	0.71
3:CD:101:VAL:HG13	3:CD:106:PHE:HD2	1.54	0.71
19:CT:19:HIS:O	19:CT:23:ARG:HG2	1.91	0.71
50:D3:4:LYS:HD2	50:D3:59:ALA:HA	1.72	0.71
23:DB:1080:A:H2'	23:DB:1081:U:C6	2.25	0.71
23:DB:1700:A:H2'	23:DB:1701:A:H5'	1.72	0.71
23:DB:455:C:N3	23:DB:472:A:H2'	2.06	0.71
28:DF:31:GLU:HG3	28:DF:158:THR:HG22	1.71	0.71
28:DF:59:ILE:O	28:DF:59:ILE:HD13	1.90	0.71
40:DS:74:ILE:HG22	40:DS:105:VAL:HG23	1.72	0.71
41:DT:49:LYS:O	41:DT:50:LEU:HG	1.90	0.71
8:AI:23:GLY:H	8:AI:60:LEU:HA	1.54	0.71
12:AM:21:ILE:HG22	12:AM:64:VAL:HG11	1.72	0.71
23:BB:1439:A:N1	23:BB:1552:A:N7	2.39	0.71
23:BB:2575:C:C5'	26:BD:149:ASN:H	2.04	0.71
25:BC:107:LYS:HB2	25:BC:194:VAL:CG1	2.20	0.71
23:BB:452:G:OP1	27:BE:52:VAL:HG22	1.90	0.71
33:BL:13:LYS:HG3	33:BL:14:LYS:N	2.01	0.71
37:BP:29:VAL:HG22	37:BP:47:ILE:HD11	1.72	0.71
38:BQ:93:ILE:HG23	38:BQ:94:LEU:H	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BT:93:LEU:HD22	41:BT:93:LEU:H	1.56	0.71
20:CB:139:GLU:O	20:CB:143:LEU:HB2	1.91	0.71
20:CB:62:ARG:H	20:CB:62:ARG:HD2	1.56	0.71
5:CF:79:ARG:HH22	5:CF:87:SER:HB3	1.56	0.71
1:CA:1186:G:H21	13:CN:100:TRP:C	1.93	0.71
50:D3:4:LYS:NZ	50:D3:60:CYS:H	1.89	0.71
23:DB:242:G:N2	23:DB:254:G:H2'	2.06	0.71
23:DB:972:A:C3'	23:DB:973:A:H5''	2.21	0.71
40:DS:17:VAL:O	40:DS:20:VAL:HG12	1.90	0.71
41:DT:29:THR:HB	41:DT:86:THR:HA	1.72	0.71
23:DB:380:G:O2'	46:DZ:13:THR:HB	1.90	0.71
1:AA:279:A:H5''	1:AA:280:C:H3'	1.72	0.71
33:BL:60:ARG:HH11	50:B3:11:LYS:HE2	1.53	0.71
23:BB:1203:U:H3'	23:BB:1204:A:H5''	1.71	0.71
23:BB:1951:U:H2'	23:BB:1953:A:OP2	1.91	0.71
23:BB:2393:U:H5'	33:BL:62:PRO:HD3	1.73	0.71
23:BB:1813:G:H21	25:BC:51:ARG:CG	2.04	0.71
1:CA:1097:C:H2'	1:CA:1098:C:C6	2.26	0.71
1:CA:1218:C:H2'	1:CA:1219:A:C8	2.26	0.71
1:CA:560:A:H5'	1:CA:566:G:N2	2.05	0.71
1:CA:859:G:H2'	1:CA:860:A:C8	2.25	0.71
23:DB:1028:A:H2'	23:DB:1029:A:C8	2.26	0.71
23:DB:1439:A:N1	23:DB:1552:A:N7	2.39	0.71
23:DB:2874:C:OP1	35:DN:5:LYS:HD2	1.91	0.71
41:DT:60:THR:HA	41:DT:82:LYS:O	1.90	0.71
45:DY:40:THR:HG23	45:DY:43:ILE:HG22	1.73	0.71
4:AE:155:LYS:HG3	7:AH:65:PHE:HB2	1.73	0.71
15:AP:20:VAL:HG23	15:AP:35:ARG:HA	1.71	0.71
47:B0:29:VAL:HB	47:B0:36:LYS:HD2	1.73	0.71
23:BB:1060:U:C4	23:BB:1088:A:N6	2.58	0.71
23:BB:1485:U:H2'	23:BB:1486:U:C6	2.26	0.71
23:BB:1700:A:H2'	23:BB:1701:A:H5'	1.72	0.71
23:BB:2039:U:H2'	23:BB:2040:G:H8	1.56	0.71
23:BB:414:C:H2'	23:BB:415:A:C8	2.25	0.71
25:BC:68:ARG:HD3	25:BC:103:ILE:HD11	1.72	0.71
29:BG:71:LEU:HA	29:BG:74:MET:SD	2.31	0.71
31:BJ:99:ARG:HA	31:BJ:103:ILE:HG13	1.73	0.71
33:BL:78:ARG:NH2	33:BL:82:LEU:HB2	2.06	0.71
36:BO:21:LEU:O	36:BO:42:PRO:HB3	1.91	0.71
36:BO:48:LEU:HD23	36:BO:49:VAL:N	2.06	0.71
3:CD:34:GLU:O	3:CD:35:GLN:HG3	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:32:THR:HG23	9:CJ:33:GLY:H	1.56	0.71
23:DB:1594:U:H2'	23:DB:1595:C:C6	2.25	0.71
23:DB:860:U:O2'	23:DB:2267:A:H4'	1.91	0.71
23:DB:950:G:H2'	23:DB:951:C:H6	1.54	0.71
29:DG:39:ALA:HB1	29:DG:54:ARG:H	1.56	0.71
30:DH:94:ILE:HG23	30:DH:98:ASP:CB	2.20	0.71
43:DW:66:VAL:HG22	43:DW:67:LYS:N	2.06	0.71
23:DB:96:C:H4'	44:DX:41:HIS:CE1	2.26	0.71
20:AB:69:VAL:O	20:AB:163:ILE:HG22	1.91	0.70
3:AD:10:LEU:HD21	3:AD:62:ARG:HD3	1.73	0.70
11:AL:87:LYS:HZ3	11:AL:87:LYS:HA	1.56	0.70
22:BA:49:C:H2'	22:BA:50:A:C8	2.26	0.70
23:BB:2037:A:H2'	23:BB:2038:G:C8	2.26	0.70
23:BB:2041:U:H2'	23:BB:2042:A:C8	2.25	0.70
23:BB:2147:A:H5'	23:BB:2148:G:C5'	2.20	0.70
23:BB:2537:U:H2'	23:BB:2538:C:C6	2.25	0.70
23:BB:279:A:H61	23:BB:361:G:H1'	1.54	0.70
25:BC:87:SER:O	25:BC:157:ALA:HB2	1.90	0.70
27:BE:151:GLY:O	27:BE:190:ALA:HB2	1.90	0.70
36:BO:16:ARG:NE	36:BO:16:ARG:HA	2.06	0.70
40:BS:28:LYS:HD3	40:BS:71:VAL:HG21	1.72	0.70
1:CA:764:C:H3'	1:CA:765:G:N2	2.06	0.70
7:CH:31:LEU:HG	7:CH:35:ILE:HD11	1.73	0.70
51:D4:34:LYS:HE2	51:D4:36:ARG:HH22	1.55	0.70
23:DB:1203:U:H3'	23:DB:1204:A:H5''	1.73	0.70
23:DB:864:G:O2'	23:DB:865:C:H5'	1.91	0.70
39:DR:79:ARG:HD3	39:DR:81:LYS:HG3	1.72	0.70
1:AA:1241:G:H2'	1:AA:1242:G:H8	1.57	0.70
1:AA:560:A:H5'	1:AA:566:G:N2	2.06	0.70
1:AA:878:A:H5''	7:AH:80:PRO:HG2	1.74	0.70
2:AC:69:THR:HG21	2:AC:75:VAL:HG21	1.73	0.70
4:AE:114:LEU:HD22	4:AE:119:VAL:HG21	1.73	0.70
50:B3:24:LYS:HZ3	50:B3:29:ARG:HH22	1.38	0.70
23:BB:1060:U:C5	52:BI:131:THR:HG22	2.25	0.70
23:BB:1469:A:H2'	23:BB:1470:A:C8	2.25	0.70
23:BB:855:G:H21	43:BW:23:LYS:CB	2.03	0.70
30:BH:90:LEU:HD21	30:BH:146:VAL:HG11	1.72	0.70
52:BI:89:SER:HB2	52:BI:136:GLY:HA3	1.73	0.70
34:BM:73:ILE:HD13	34:BM:92:TRP:H	1.56	0.70
41:BT:61:LEU:HD12	41:BT:62:VAL:H	1.56	0.70
3:CD:187:ARG:HH12	3:CD:191:SER:HA	1.55	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CE:87:VAL:HG22	4:CE:88:HIS:H	1.55	0.70
13:CN:51:PRO:HG2	13:CN:52:ARG:H	1.56	0.70
50:D3:51:LYS:HA	50:D3:51:LYS:HZ2	1.56	0.70
23:DB:2306:C:H3'	23:DB:2307:G:H5''	1.72	0.70
23:DB:28:A:N6	23:DB:512:G:O2'	2.24	0.70
23:DB:350:G:H2'	23:DB:351:C:O4'	1.91	0.70
31:DJ:35:ARG:HB3	31:DJ:54:ILE:HD11	1.72	0.70
42:DU:43:LYS:HG2	42:DU:57:ILE:HB	1.73	0.70
5:AF:53:LYS:HD3	5:AF:54:LEU:N	2.04	0.70
18:AS:4:LEU:C	18:AS:5:LYS:HE3	2.12	0.70
23:BB:1182:G:H2'	23:BB:1183:U:O4'	1.90	0.70
23:BB:131:A:H2'	23:BB:132:G:H8	1.57	0.70
23:BB:2197:U:O2'	23:BB:2198:A:H2'	1.91	0.70
25:BC:10:PRO:HD2	25:BC:202:ARG:HH12	1.57	0.70
23:BB:2483:C:H4'	34:BM:51:ARG:HH12	1.56	0.70
1:CA:1238:A:H5'	1:CA:1336:C:N4	2.04	0.70
1:CA:451:A:H4'	1:CA:452:A:O4'	1.91	0.70
12:CM:52:ILE:HA	12:CM:55:LEU:HD12	1.72	0.70
48:D1:40:PRO:HD2	48:D1:44:GLN:O	1.90	0.70
51:D4:30:GLU:HB3	51:D4:33:HIS:HB2	1.72	0.70
23:DB:1951:U:H2'	23:DB:1953:A:OP2	1.91	0.70
23:DB:2037:A:H2'	23:DB:2038:G:C8	2.26	0.70
23:DB:2557:G:H2'	23:DB:2558:C:C6	2.27	0.70
25:DC:164:VAL:HG23	25:DC:167:ASP:HB2	1.71	0.70
26:DD:61:THR:HG23	26:DD:62:LYS:H	1.54	0.70
43:DW:42:THR:H	43:DW:65:LYS:HG2	1.57	0.70
4:AE:101:GLY:H	4:AE:121:ASN:HD21	1.39	0.70
23:BB:1082:U:C4	23:BB:1086:A:C2	2.80	0.70
23:BB:2228:G:H2'	23:BB:2229:U:C6	2.26	0.70
23:BB:287:G:H2'	23:BB:288:U:H6	1.55	0.70
26:BD:154:LYS:C	26:BD:156:PHE:H	1.94	0.70
29:BG:19:ASN:HB3	29:BG:22:VAL:HG13	1.72	0.70
29:BG:40:VAL:HB	29:BG:63:GLN:HE22	1.57	0.70
31:BJ:41:LYS:HD2	31:BJ:44:TYR:HB2	1.73	0.70
34:BM:33:LEU:N	34:BM:101:VAL:HG22	2.06	0.70
34:BM:70:ASP:H	34:BM:71:LYS:NZ	1.90	0.70
35:BN:24:MET:HE3	35:BN:44:LEU:HD22	1.74	0.70
46:BZ:15:SER:HB2	46:BZ:25:ARG:NH1	2.06	0.70
23:DB:2537:U:H2'	23:DB:2538:C:C6	2.26	0.70
23:DB:98:G:H3'	23:DB:99:U:H5''	1.71	0.70
25:DC:28:PRO:HG2	25:DC:79:ARG:NE	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:108:ASP:HB2	26:DD:204:LYS:HB2	1.74	0.70
27:DE:143:LEU:HD22	27:DE:143:LEU:N	2.06	0.70
28:DF:102:LEU:O	28:DF:107:VAL:HB	1.91	0.70
23:DB:1190:G:OP1	33:DL:39:LYS:N	2.24	0.70
37:DP:76:HIS:CD2	37:DP:76:HIS:N	2.60	0.70
42:DU:71:ILE:HD12	42:DU:102:ILE:HD12	1.74	0.70
2:AC:5:HIS:HD2	2:AC:8:GLY:H	1.39	0.70
16:AQ:76:ARG:NE	16:AQ:78:VAL:HG22	2.04	0.70
50:B3:29:ARG:HA	50:B3:33:THR:CG2	2.22	0.70
23:BB:990:A:H1'	23:BB:1156:A:C2	2.26	0.70
23:BB:1594:U:H2'	23:BB:1595:C:C6	2.26	0.70
23:BB:191:A:H2'	23:BB:192:C:C6	2.26	0.70
40:BS:18:ARG:HH12	40:BS:25:ARG:HH22	1.38	0.70
41:BT:62:VAL:O	41:BT:63:VAL:HG12	1.90	0.70
23:BB:85:G:C5'	42:BU:28:LEU:HA	2.17	0.70
24:BV:80:HIS:HD2	24:BV:83:LYS:H	1.39	0.70
1:CA:865:A:H5'	1:CA:1078:U:O4	1.92	0.70
1:CA:545:C:O2'	1:CA:546:A:H5'	1.90	0.70
10:CK:28:ASN:HD21	10:CK:47:GLY:H	1.39	0.70
21:CU:36:PHE:HB2	21:CU:39:LYS:HB2	1.74	0.70
47:D0:10:SER:O	47:D0:14:MET:HB2	1.91	0.70
22:DA:61:G:H2'	22:DA:62:C:H6	1.55	0.70
23:DB:1794:A:H2'	23:DB:1795:C:H6	1.57	0.70
25:DC:29:PHE:HD1	25:DC:29:PHE:O	1.73	0.70
37:DP:23:ASP:HA	37:DP:49:ILE:HG22	1.73	0.70
13:AN:74:ARG:HD3	13:AN:74:ARG:O	1.90	0.70
23:BB:1826:G:OP2	25:BC:221:GLY:HA2	1.90	0.70
23:BB:2216:G:H2'	23:BB:2217:G:C8	2.26	0.70
25:BC:172:THR:HG22	25:BC:173:LEU:H	1.54	0.70
29:BG:28:LYS:HD3	29:BG:79:THR:HA	1.72	0.70
3:CD:94:GLU:HG2	3:CD:185:PRO:HG3	1.73	0.70
9:CJ:11:LYS:HB2	9:CJ:97:ASP:HB3	1.74	0.70
10:CK:16:SER:CA	10:CK:78:ILE:HA	2.20	0.70
23:DB:596:U:H2'	23:DB:597:G:C8	2.26	0.70
30:DH:50:ARG:HG3	30:DH:51:ARG:N	2.06	0.70
31:DJ:60:ASP:HB3	31:DJ:126:ALA:HB1	1.73	0.70
23:BB:1015:U:H2'	23:BB:1016:G:H8	1.57	0.70
23:BB:181:A:H2'	23:BB:182:A:H8	1.56	0.70
14:AO:88:ARG:HG3	23:BB:716:A:OP1	1.91	0.70
27:BE:106:LYS:HB2	27:BE:106:LYS:NZ	2.07	0.70
39:BR:18:GLN:HB3	39:BR:99:THR:HA	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BV:80:HIS:HD2	24:BV:82:TYR:H	1.37	0.70
43:BW:17:ALA:HA	43:BW:35:ILE:O	1.92	0.70
1:CA:17:U:H2'	1:CA:18:C:H6	1.56	0.70
47:D0:26:SER:HB2	47:D0:38:LEU:HD21	1.74	0.70
23:DB:1099:G:H4'	52:DI:4:VAL:HB	1.73	0.70
23:DB:1178:C:H2'	23:DB:1179:G:C8	2.27	0.70
23:DB:570:G:H2'	23:DB:2030:A:N7	2.05	0.70
23:DB:2246:G:H2'	23:DB:2247:A:C8	2.27	0.70
23:DB:2684:U:OP1	37:DP:59:THR:HB	1.91	0.70
25:DC:141:HIS:CB	25:DC:190:THR:HB	2.22	0.70
35:DN:2:ARG:HH11	35:DN:2:ARG:HG2	1.57	0.70
36:DO:35:ILE:HG12	36:DO:106:LEU:HD12	1.74	0.70
36:DO:38:GLN:HA	36:DO:50:ALA:HB3	1.72	0.70
1:AA:1219:A:H2'	1:AA:1220:G:C8	2.27	0.70
1:AA:1238:A:H5'	1:AA:1336:C:H41	1.55	0.70
23:BB:929:U:H5'	45:BY:37:ARG:NH1	2.07	0.70
30:BH:17:ASP:OD2	30:BH:21:VAL:HB	1.92	0.70
38:BQ:92:LYS:CG	38:BQ:93:ILE:N	2.54	0.70
38:BQ:92:LYS:HA	38:BQ:95:ALA:HB2	1.72	0.70
5:CF:38:ARG:HD3	5:CF:97:THR:HA	1.73	0.70
25:DC:144:GLU:HB2	25:DC:187:CYS:HB2	1.73	0.70
23:DB:2597:G:H5''	25:DC:239:PHE:HB2	1.74	0.70
30:DH:8:LYS:HE2	30:DH:9:VAL:H	1.56	0.70
39:DR:70:GLU:HG3	39:DR:71:LYS:HG2	1.73	0.70
46:DZ:5:ILE:HG13	46:DZ:51:VAL:HG13	1.73	0.70
1:AA:1289:A:H2'	1:AA:1290:G:H5'	1.74	0.70
8:AI:40:ARG:N	8:AI:44:ARG:HE	1.88	0.70
17:AR:58:ILE:O	17:AR:62:ARG:HG3	1.92	0.70
23:BB:1794:A:H2'	23:BB:1795:C:C6	2.26	0.70
23:BB:1794:A:H2'	23:BB:1795:C:H6	1.57	0.70
23:BB:2297:A:H61	23:BB:2319:G:H3'	1.54	0.70
23:BB:2834:G:H1'	23:BB:2883:A:N6	2.06	0.70
34:BM:23:GLY:N	34:BM:96:ILE:HD12	2.06	0.70
22:BA:98:G:H1	24:BV:14:LYS:HB2	1.55	0.70
46:BZ:69:SER:O	46:BZ:70:LYS:HB2	1.91	0.70
1:CA:1007:U:H2'	1:CA:1008:U:C6	2.27	0.70
1:CA:1216:A:OP1	13:CN:2:LYS:HE2	1.91	0.70
1:CA:1391:U:H2'	1:CA:1392:G:C8	2.26	0.70
1:CA:188:C:H2'	1:CA:189:A:O4'	1.90	0.70
1:CA:960:U:O2'	1:CA:961:U:OP2	2.09	0.70
1:CA:1342:C:H5'	8:CI:127:SER:HB3	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CS:48:ILE:O	18:CS:59:VAL:HG22	1.90	0.70
23:DB:1113:U:H5'	29:DG:2:ARG:HG2	1.72	0.70
23:DB:1794:A:H2'	23:DB:1795:C:C6	2.26	0.70
39:DR:96:VAL:HG12	39:DR:98:ILE:HG12	1.72	0.70
41:DT:17:SER:H	41:DT:20:ALA:HB3	1.55	0.70
1:AA:1071:C:H2'	1:AA:1072:G:H8	1.55	0.70
3:AD:2:ARG:NH1	3:AD:114:ARG:HD2	2.05	0.70
3:AD:25:ARG:HD3	3:AD:26:ALA:H	1.56	0.70
23:BB:2557:G:H2'	23:BB:2558:C:C6	2.27	0.70
42:BU:25:LYS:HG2	42:BU:35:VAL:HG22	1.74	0.70
24:BV:40:ILE:H	24:BV:40:ILE:HD13	1.56	0.70
43:BW:36:ILE:HB	43:BW:68:PHE:HD2	1.57	0.70
1:CA:939:G:H5''	6:CG:101:ARG:NH1	2.06	0.70
23:DB:1064:C:C4'	52:DI:90:GLY:HA2	2.21	0.70
23:DB:1485:U:H2'	23:DB:1486:U:C6	2.26	0.70
23:DB:28:A:H61	23:DB:512:G:H1'	1.57	0.70
23:DB:969:G:H2'	23:DB:970:U:C6	2.26	0.70
31:DJ:84:ILE:CD1	31:DJ:85:LYS:H	2.04	0.70
32:DK:110:GLU:HA	32:DK:113:MET:HG2	1.74	0.70
41:DT:68:LYS:HB2	41:DT:68:LYS:NZ	2.07	0.70
14:AO:11:VAL:HG23	14:AO:26:VAL:HG11	1.72	0.69
16:AQ:10:ARG:HH11	16:AQ:55:GLY:H	1.37	0.69
23:BB:150:U:H2'	23:BB:151:C:C6	2.27	0.69
23:BB:2270:A:H4'	43:BW:18:LYS:CG	2.21	0.69
26:BD:119:ALA:HA	26:BD:123:LYS:NZ	2.06	0.69
23:BB:1188:U:H4'	39:BR:83:TYR:HB2	1.74	0.69
20:CB:8:MET:SD	20:CB:9:LEU:HD22	2.32	0.69
23:DB:1346:G:O2'	23:DB:1347:A:H5'	1.92	0.69
25:DC:63:ILE:HG21	25:DC:90:ILE:CD1	2.22	0.69
23:DB:336:C:H5''	42:DU:3:LYS:NZ	2.06	0.69
1:AA:1236:A:H4'	1:AA:1304:G:H4'	1.73	0.69
14:AO:87:ARG:HA	14:AO:87:ARG:NH1	2.07	0.69
23:BB:2619:C:H5'	26:BD:157:LYS:CE	2.21	0.69
35:BN:33:ILE:HD13	35:BN:33:ILE:H	1.56	0.69
35:BN:4:ARG:HD2	35:BN:5:LYS:H	1.56	0.69
37:BP:32:VAL:HG13	37:BP:40:GLN:HE21	1.57	0.69
43:BW:10:ARG:HB3	43:BW:10:ARG:HH11	1.57	0.69
1:CA:279:A:H5''	1:CA:280:C:H3'	1.72	0.69
7:CH:55:LYS:NZ	7:CH:55:LYS:HA	2.07	0.69
16:CQ:19:SER:O	16:CQ:20:ILE:HD12	1.93	0.69
23:DB:172:A:H2'	23:DB:173:A:C8	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2147:A:H4'	23:DB:2148:G:H8	1.57	0.69
23:DB:522:A:H2'	23:DB:523:C:C6	2.27	0.69
23:DB:920:A:H2'	23:DB:921:C:C6	2.27	0.69
23:DB:2636:C:H5'	26:DD:80:TRP:HZ2	1.57	0.69
29:DG:155:PRO:CB	29:DG:171:LYS:HB3	2.21	0.69
37:DP:31:VAL:HG13	37:DP:81:ASP:HB3	1.73	0.69
40:DS:73:LYS:HD3	40:DS:74:ILE:N	2.06	0.69
46:DZ:47:LYS:HB2	46:DZ:51:VAL:HG12	1.74	0.69
1:AA:473:U:H2'	1:AA:474:G:H8	1.58	0.69
23:BB:1373:A:H2'	23:BB:1374:G:O4'	1.92	0.69
23:BB:1406:U:H2'	23:BB:1407:G:C8	2.27	0.69
23:BB:150:U:H2'	23:BB:151:C:H6	1.56	0.69
23:BB:38:A:H2'	23:BB:39:G:O4'	1.92	0.69
26:BD:11:MET:SD	37:BP:9:GLN:HG2	2.32	0.69
27:BE:146:VAL:O	27:BE:184:ASP:HB2	1.92	0.69
34:BM:133:LYS:O	34:BM:133:LYS:HE3	1.93	0.69
1:CA:1320:C:N3	18:CS:35:ARG:HD3	2.06	0.69
2:CC:39:ARG:HH22	13:CN:91:GLU:HB3	1.58	0.69
51:D4:1:MET:HB3	51:D4:2:LYS:HD3	1.74	0.69
23:DB:1098:A:H2'	52:DI:4:VAL:CA	2.22	0.69
23:DB:151:C:H2'	23:DB:152:A:C8	2.27	0.69
28:DF:162:ASP:O	28:DF:166:ARG:HG3	1.92	0.69
1:AA:413:G:O6	3:AD:32:LYS:HE3	1.93	0.69
20:AB:59:ILE:HD12	20:AB:60:ALA:N	2.07	0.69
21:AU:36:PHE:HB3	21:AU:40:PRO:CD	2.21	0.69
51:B4:24:ARG:HA	51:B4:34:LYS:O	1.93	0.69
25:BC:164:VAL:HB	25:BC:167:ASP:OD1	1.92	0.69
31:BJ:68:LYS:HA	31:BJ:71:ASP:OD1	1.93	0.69
36:BO:41:ALA:HB1	36:BO:42:PRO:HD2	1.74	0.69
37:BP:18:SER:HA	37:BP:87:ARG:NH2	2.07	0.69
37:BP:2:ASN:N	37:BP:2:ASN:HD22	1.89	0.69
40:BS:7:HIS:HB2	40:BS:103:ILE:HG22	1.72	0.69
41:BT:64:LYS:HD2	41:BT:77:ARG:HD3	1.75	0.69
42:BU:46:LYS:N	42:BU:47:PRO:HD3	2.08	0.69
1:CA:1508:A:H2'	1:CA:1509:C:H6	1.56	0.69
12:CM:92:ARG:CZ	12:CM:92:ARG:HA	2.22	0.69
23:DB:2578:G:O2'	26:DD:138:LEU:HD13	1.93	0.69
23:DB:2834:G:H1'	23:DB:2883:A:N6	2.06	0.69
26:DD:30:GLU:HG2	26:DD:94:GLN:NE2	2.07	0.69
23:DB:2636:C:H5'	26:DD:80:TRP:CZ2	2.28	0.69
27:DE:83:VAL:O	27:DE:84:THR:HG22	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:66:LYS:N	41:DT:76:ARG:HH21	1.87	0.69
24:DV:21:ARG:HE	24:DV:87:GLN:HA	1.57	0.69
1:AA:1432:G:OP1	37:BP:105:LYS:HB2	1.93	0.69
22:BA:6:G:H2'	22:BA:7:G:C8	2.27	0.69
23:BB:2469:A:H4'	34:BM:55:ARG:HG3	1.75	0.69
23:BB:45:G:C5'	23:BB:46:G:H5'	2.22	0.69
37:BP:77:SER:H	37:BP:78:PRO:CD	2.06	0.69
24:BV:77:VAL:HA	24:BV:89:ILE:HG22	1.74	0.69
11:CL:51:VAL:HG12	11:CL:52:CYS:H	1.57	0.69
48:D1:45:HIS:O	48:D1:46:VAL:HG12	1.92	0.69
25:DC:53:ILE:HG12	25:DC:218:THR:HA	1.74	0.69
27:DE:136:GLN:O	27:DE:139:LYS:HG2	1.93	0.69
27:DE:163:ASN:H	27:DE:168:ASP:HA	1.58	0.69
41:DT:56:GLU:HG2	41:DT:57:VAL:HG13	1.75	0.69
1:AA:1048:G:H4'	13:AN:2:LYS:NZ	2.07	0.69
1:AA:270:A:H2'	1:AA:271:C:H6	1.57	0.69
1:AA:60:A:H4'	1:AA:61:G:O5'	1.92	0.69
20:AB:76:SER:O	20:AB:79:VAL:HG12	1.93	0.69
15:AP:5:ARG:HD2	15:AP:5:ARG:H	1.57	0.69
16:AQ:11:VAL:HA	16:AQ:22:VAL:HG22	1.73	0.69
5:AF:86:ARG:HH11	17:AR:63:TYR:HB3	1.57	0.69
47:B0:30:ASP:OD1	47:B0:31:LYS:HD2	1.93	0.69
50:B3:26:ALA:HB1	50:B3:29:ARG:HG2	1.73	0.69
23:BB:2037:A:H2'	23:BB:2038:G:H8	1.58	0.69
23:BB:2241:A:H2'	23:BB:2242:G:H8	1.57	0.69
23:BB:2645:G:H3'	23:BB:2646:C:C5'	2.23	0.69
23:BB:2804:U:H2'	23:BB:2805:C:C6	2.28	0.69
23:BB:596:U:H2'	23:BB:597:G:C8	2.27	0.69
25:BC:170:TYR:O	25:BC:171:VAL:HG13	1.91	0.69
27:BE:98:LYS:HG3	27:BE:99:LYS:H	1.56	0.69
31:BJ:6:ALA:O	31:BJ:8:PRO:HD3	1.93	0.69
32:BK:85:VAL:HG11	32:BK:115:ILE:HD12	1.74	0.69
34:BM:69:PRO:HB2	34:BM:71:LYS:NZ	2.08	0.69
38:BQ:102:LYS:H	38:BQ:102:LYS:HD2	1.58	0.69
38:BQ:96:ASP:O	38:BQ:99:VAL:HG22	1.92	0.69
39:BR:79:ARG:HD3	39:BR:86:GLN:NE2	2.08	0.69
1:CA:1250:A:H2'	1:CA:1251:A:C8	2.28	0.69
1:CA:832:G:O2'	1:CA:833:G:H5'	1.92	0.69
10:CK:85:VAL:HG21	10:CK:92:ARG:HH12	1.58	0.69
23:DB:2017:U:O2	47:D0:6:LYS:HG3	1.92	0.69
23:DB:1309:G:H5'	49:D2:8:SER:H	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1552:A:H2'	23:DB:1553:A:H5'	1.73	0.69
1:AA:1218:C:H2'	1:AA:1219:A:C8	2.28	0.69
1:AA:978:A:H5'	1:AA:1362:A:N6	2.08	0.69
51:B4:24:ARG:HH22	51:B4:36:ARG:CD	2.03	0.69
25:BC:141:HIS:CD2	25:BC:193:GLU:H	2.11	0.69
2:CC:11:LEU:HD22	2:CC:17:TRP:HE1	1.58	0.69
8:CI:118:ARG:NH2	8:CI:122:ARG:HE	1.90	0.69
22:DA:66:A:O2'	22:DA:67:G:H5''	1.93	0.69
23:DB:1118:C:H2'	23:DB:1119:U:C6	2.27	0.69
23:DB:1866:A:H2'	23:DB:1867:G:O4'	1.93	0.69
23:DB:2229:U:H2'	23:DB:2230:G:C8	2.27	0.69
23:DB:2645:G:H3'	23:DB:2646:C:C5'	2.22	0.69
27:DE:15:SER:HB3	27:DE:196:VAL:HG22	1.74	0.69
27:DE:46:GLN:NE2	27:DE:86:ALA:HB3	2.07	0.69
30:DH:5:LEU:HD13	30:DH:35:LYS:HE3	1.74	0.69
31:DJ:135:GLN:HE21	31:DJ:138:GLN:H	1.40	0.69
36:DO:15:ARG:HH12	43:DW:76:ARG:NE	1.91	0.69
1:AA:451:A:H4'	1:AA:452:A:O4'	1.91	0.69
23:BB:1799:G:H4'	23:BB:1800:C:O5'	1.93	0.69
23:BB:437:U:H2'	23:BB:438:G:C8	2.27	0.69
25:BC:127:ASN:HD22	25:BC:128:THR:N	1.91	0.69
23:BB:1243:C:O2'	33:BL:15:ALA:HB3	1.92	0.69
1:CA:1281:C:H5'	1:CA:1282:C:H5	1.56	0.69
1:CA:412:A:H1'	1:CA:413:G:H8	1.57	0.69
48:D1:29:LYS:HB2	48:D1:30:PRO:CD	2.18	0.69
23:DB:1098:A:OP2	52:DI:3:LYS:HG2	1.92	0.69
23:DB:682:G:H5'	49:D2:26:ASN:ND2	2.08	0.69
31:DJ:124:VAL:HG23	31:DJ:125:TYR:N	2.07	0.69
39:DR:22:LEU:HD12	39:DR:23:GLU:H	1.57	0.69
42:DU:91:LYS:HD3	42:DU:93:ARG:NE	2.07	0.69
1:AA:1060:U:H2'	1:AA:1061:G:H8	1.58	0.69
20:AB:53:LEU:HA	20:AB:56:LEU:HD22	1.74	0.69
8:AI:18:VAL:HG11	8:AI:82:ILE:HG12	1.75	0.69
23:BB:1252:G:H1	38:BQ:36:GLN:HE22	1.39	0.69
23:BB:215:G:H4'	23:BB:216:A:H4'	1.75	0.69
25:BC:79:ARG:HH11	25:BC:80:LEU:H	1.38	0.69
30:BH:82:SER:N	30:BH:146:VAL:HG13	2.00	0.69
31:BJ:28:LEU:HD23	31:BJ:29:ALA:N	2.08	0.69
23:BB:2355:G:H4'	43:BW:20:LEU:HD22	1.75	0.69
1:CA:1096:C:H2'	1:CA:1097:C:C6	2.28	0.69
1:CA:473:U:H2'	1:CA:474:G:H8	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:39:PRO:HA	9:CJ:74:VAL:HG22	1.74	0.69
22:DA:22:U:H2'	22:DA:23:G:C8	2.28	0.69
23:DB:2834:G:H1'	23:DB:2883:A:H61	1.58	0.69
27:DE:108:ILE:HG13	27:DE:109:LEU:H	1.58	0.69
32:DK:64:ARG:HD2	32:DK:102:PRO:O	1.91	0.69
39:DR:3:ALA:O	39:DR:4:VAL:HG13	1.93	0.69
39:DR:49:ILE:HG13	39:DR:51:VAL:HG23	1.74	0.69
23:DB:993:G:H21	39:DR:93:PHE:HZ	1.41	0.69
45:DY:7:THR:HA	45:DY:34:THR:HB	1.75	0.69
46:DZ:25:ARG:HG3	46:DZ:26:SER:H	1.58	0.69
1:AA:131:A:H2'	1:AA:132:C:C6	2.28	0.69
1:AA:1524:C:H2'	1:AA:1525:G:C8	2.28	0.69
1:AA:832:G:O2'	1:AA:833:G:H5'	1.92	0.69
2:AC:9:ILE:HG23	2:AC:10:ARG:HG3	1.73	0.69
5:AF:3:HIS:HA	5:AF:65:GLU:HA	1.75	0.69
7:AH:87:ARG:HB2	7:AH:90:GLU:HG3	1.74	0.69
10:AK:109:ILE:H	21:AU:5:VAL:HB	1.57	0.69
47:B0:26:SER:HB2	47:B0:38:LEU:HD12	1.74	0.69
47:B0:29:VAL:HG21	47:B0:41:HIS:ND1	2.08	0.69
23:BB:1866:A:H2'	23:BB:1867:G:O4'	1.93	0.69
23:BB:743:A:O2'	23:BB:744:U:H5'	1.93	0.69
29:BG:97:VAL:HG11	29:BG:122:ALA:O	1.92	0.69
34:BM:100:LYS:O	34:BM:100:LYS:HD2	1.93	0.69
42:BU:90:LYS:HZ2	42:BU:91:LYS:HE2	1.57	0.69
49:D2:21:ARG:HG3	49:D2:31:LEU:HD11	1.75	0.69
27:DE:115:GLN:HG3	27:DE:116:ASP:H	1.58	0.69
29:DG:95:ALA:HB1	29:DG:130:ILE:HD11	1.73	0.69
52:DI:85:ILE:HD13	52:DI:137:LEU:HD21	1.75	0.69
23:DB:1099:G:O5'	52:DI:4:VAL:HG12	1.92	0.69
33:DL:64:PHE:HA	50:D3:11:LYS:HA	1.75	0.69
34:DM:5:LYS:HG3	34:DM:69:PRO:O	1.93	0.69
36:DO:7:ARG:O	36:DO:11:ALA:HB2	1.93	0.69
37:DP:6:GLN:CD	37:DP:6:GLN:H	1.95	0.69
40:DS:48:LYS:O	40:DS:52:GLU:HG3	1.92	0.69
46:DZ:3:LYS:HG3	46:DZ:48:GLN:H	1.58	0.69
8:AI:12:LYS:H	8:AI:109:GLN:HE22	1.41	0.69
50:B3:31:ILE:HG22	50:B3:32:LEU:HG	1.75	0.69
23:BB:2265:U:H3'	23:BB:2266:A:H5''	1.75	0.69
23:BB:845:A:C2	23:BB:847:U:H1'	2.27	0.69
25:BC:68:ARG:HD2	25:BC:127:ASN:HD21	1.57	0.69
42:BU:54:PRO:HG2	42:BU:55:GLY:H	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1218:C:H2'	1:CA:1219:A:H8	1.56	0.69
1:CA:131:A:H2'	1:CA:132:C:C6	2.28	0.69
23:DB:1116:G:H21	34:DM:136:MET:HE3	1.57	0.69
23:DB:1440:U:H2'	23:DB:1441:G:C8	2.28	0.69
23:DB:2229:U:H2'	23:DB:2230:G:H8	1.56	0.69
34:DM:64:TRP:HB2	34:DM:102:LEU:HB2	1.75	0.69
40:DS:83:LYS:HD3	40:DS:83:LYS:N	2.07	0.69
1:AA:865:A:H5'	1:AA:1078:U:O4	1.93	0.68
1:AA:1330:U:C2'	1:AA:1331:G:H5'	2.22	0.68
1:AA:412:A:H1'	1:AA:413:G:H8	1.57	0.68
1:AA:673:A:H2'	1:AA:674:G:C8	2.29	0.68
3:AD:60:VAL:HA	3:AD:63:ILE:HD12	1.74	0.68
23:BB:1060:U:O2	23:BB:1088:A:N7	2.27	0.68
23:BB:1791:A:H1'	25:BC:206:LYS:NZ	2.08	0.68
23:BB:2849:U:H4'	23:BB:2850:A:H5'	1.74	0.68
33:BL:118:THR:CB	33:BL:119:PRO:HD3	2.22	0.68
36:BO:68:LYS:HD3	36:BO:68:LYS:H	1.58	0.68
40:BS:7:HIS:CD2	40:BS:8:ARG:H	2.11	0.68
1:CA:270:A:H2'	1:CA:271:C:H6	1.57	0.68
1:CA:642:A:H2'	1:CA:643:C:H6	1.58	0.68
5:CF:47:LEU:HD12	5:CF:55:HIS:HA	1.75	0.68
48:D1:10:LEU:H	48:D1:24:LYS:HB3	1.58	0.68
23:DB:532:A:N1	23:DB:2020:A:H1'	2.08	0.68
33:DL:35:HIS:NE2	39:DR:84:ARG:HG2	2.08	0.68
39:DR:37:GLU:HG3	39:DR:62:GLU:HG3	1.74	0.68
4:AE:155:LYS:HE3	7:AH:70:VAL:HG13	1.75	0.68
9:AJ:50:THR:HG22	9:AJ:64:GLN:HG2	1.75	0.68
18:AS:20:LYS:HB3	18:AS:20:LYS:HZ2	1.56	0.68
23:BB:1346:G:O2'	23:BB:1347:A:H5'	1.92	0.68
23:BB:1552:A:H2'	23:BB:1553:A:H5'	1.74	0.68
23:BB:181:A:H2'	23:BB:182:A:C8	2.28	0.68
23:BB:423:A:H5'	23:BB:424:G:H5''	1.75	0.68
25:BC:124:LYS:HB2	25:BC:125:PRO:HD3	1.75	0.68
29:BG:120:ILE:HG12	29:BG:140:ILE:HG22	1.74	0.68
31:BJ:119:PHE:CG	31:BJ:120:ARG:N	2.61	0.68
31:BJ:36:LEU:HD11	31:BJ:122:LEU:HD23	1.75	0.68
37:BP:59:THR:O	37:BP:60:VAL:HG13	1.94	0.68
39:BR:41:ILE:C	39:BR:43:ASN:H	1.95	0.68
46:BZ:5:ILE:HB	46:BZ:51:VAL:CG1	2.23	0.68
1:CA:203:G:H1'	1:CA:465:A:H62	1.58	0.68
1:CA:977:A:H3'	1:CA:1362:A:H62	1.56	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:98:A:H2'	1:CA:99:C:C6	2.28	0.68
20:CB:83:ALA:O	20:CB:88:GLN:HB2	1.93	0.68
2:CC:156:LEU:HB2	2:CC:163:ARG:HD3	1.73	0.68
8:CI:20:ILE:HG23	8:CI:60:LEU:HD13	1.74	0.68
21:CU:14:ALA:H	21:CU:16:ARG:HH22	1.41	0.68
23:DB:1406:U:H2'	23:DB:1407:G:C8	2.28	0.68
23:DB:2054:A:H2'	47:D0:4:GLN:NE2	2.08	0.68
32:DK:64:ARG:H	32:DK:83:ALA:HB3	1.58	0.68
23:BB:2071:A:H2'	23:BB:2072:C:C6	2.28	0.68
23:BB:2901:C:H2'	23:BB:2902:C:H5'	1.74	0.68
23:BB:974:G:H1'	23:BB:975:A:C8	2.27	0.68
25:BC:61:TYR:HD2	25:BC:84:PRO:HD2	1.58	0.68
28:BF:46:LYS:O	28:BF:50:ASP:HB3	1.94	0.68
29:BG:176:LYS:O	29:BG:176:LYS:HE2	1.93	0.68
31:BJ:93:ILE:HD13	31:BJ:100:VAL:HG21	1.74	0.68
23:BB:65:U:H4'	41:BT:70:HIS:ND1	2.09	0.68
1:CA:1328:C:H5''	12:CM:27:THR:HG21	1.75	0.68
23:DB:2037:A:H2'	23:DB:2038:G:H8	1.58	0.68
23:DB:2246:G:H2'	23:DB:2247:A:H8	1.58	0.68
23:DB:2804:U:H2'	23:DB:2805:C:C6	2.27	0.68
23:DB:2804:U:H2'	23:DB:2805:C:H6	1.58	0.68
23:DB:2860:A:H8	23:DB:2860:A:O5'	1.76	0.68
25:DC:89:ASN:HB2	25:DC:105:ALA:HB3	1.75	0.68
29:DG:17:LYS:HE2	29:DG:24:THR:H	1.59	0.68
43:DW:56:HIS:HA	43:DW:77:LYS:HE2	1.75	0.68
46:DZ:39:LYS:HD3	46:DZ:61:ASN:ND2	2.09	0.68
20:AB:9:LEU:HD22	20:AB:11:ALA:N	2.07	0.68
23:BB:2795:C:H2'	23:BB:2796:U:O4'	1.94	0.68
23:BB:2860:A:O5'	23:BB:2860:A:H8	1.76	0.68
23:BB:594:U:H2'	23:BB:595:C:C6	2.28	0.68
23:BB:720:U:H2'	23:BB:721:A:C8	2.28	0.68
23:BB:1803:A:O3'	25:BC:254:LYS:HG2	1.94	0.68
33:BL:30:THR:HG22	33:BL:36:LYS:NZ	2.07	0.68
42:BU:25:LYS:HA	42:BU:35:VAL:H	1.58	0.68
43:BW:43:LYS:HG2	43:BW:76:ARG:HE	1.57	0.68
1:CA:1057:G:H4'	2:CC:196:GLY:H	1.58	0.68
3:CD:141:VAL:HA	3:CD:180:THR:H	1.58	0.68
8:CI:70:GLY:O	8:CI:74:GLN:HG2	1.93	0.68
51:D4:15:LYS:NZ	51:D4:22:VAL:HG12	2.06	0.68
51:D4:26:ILE:HG13	51:D4:35:GLN:N	2.03	0.68
23:DB:1082:U:C4	23:DB:1086:A:C2	2.81	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:947:A:H2'	23:DB:948:C:C6	2.29	0.68
26:DD:79:LEU:HG	26:DD:80:TRP:H	1.57	0.68
30:DH:26:ALA:HB3	30:DH:31:VAL:HG23	1.74	0.68
1:AA:642:A:H2'	1:AA:643:C:H6	1.58	0.68
9:AJ:56:HIS:H	13:AN:80:ARG:HH22	1.41	0.68
10:AK:70:ALA:HB1	10:AK:74:LYS:HE3	1.75	0.68
10:AK:97:ARG:NH1	10:AK:97:ARG:HB3	2.08	0.68
22:BA:17:C:O2'	22:BA:18:G:H5'	1.93	0.68
23:BB:139:U:H5''	23:BB:140:C:C5	2.27	0.68
23:BB:1440:U:H2'	23:BB:1441:G:C8	2.28	0.68
27:BE:113:VAL:HG23	27:BE:117:ARG:NH2	2.09	0.68
27:BE:156:ASN:ND2	27:BE:157:LEU:HG	2.09	0.68
23:BB:322:A:OP1	27:BE:162:ARG:HG2	1.93	0.68
39:BR:39:LEU:O	39:BR:54:VAL:CG1	2.41	0.68
44:BX:30:MET:HG3	44:BX:31:GLN:H	1.58	0.68
18:CS:10:ILE:HB	18:CS:14:LEU:HD11	1.75	0.68
18:CS:51:HIS:HA	18:CS:56:HIS:HA	1.75	0.68
50:D3:32:LEU:CD1	50:D3:33:THR:H	2.06	0.68
23:DB:1098:A:C2'	52:DI:4:VAL:N	2.55	0.68
23:DB:533:G:H2'	23:DB:534:U:C6	2.28	0.68
23:DB:857:G:O2'	23:DB:858:G:H5'	1.94	0.68
25:DC:63:ILE:HD13	25:DC:90:ILE:HD13	1.76	0.68
27:DE:161:ALA:O	27:DE:169:VAL:HB	1.93	0.68
31:DJ:19:ASP:HB3	31:DJ:21:THR:HG23	1.75	0.68
33:DL:124:GLY:H	33:DL:142:ILE:CA	2.07	0.68
35:DN:97:ILE:HA	35:DN:113:ILE:HD13	1.74	0.68
38:DQ:108:LEU:HA	38:DQ:111:LYS:HD2	1.76	0.68
1:AA:1342:C:H2'	1:AA:1343:G:C8	2.28	0.68
1:AA:194:C:O2'	1:AA:195:A:H5'	1.94	0.68
11:AL:8:ARG:CZ	11:AL:9:LYS:HE3	2.24	0.68
12:AM:77:LYS:HA	12:AM:80:MET:HB3	1.74	0.68
15:AP:68:SER:HB3	15:AP:71:VAL:HG12	1.76	0.68
23:BB:303:G:H2'	23:BB:304:U:C6	2.29	0.68
25:BC:20:ASN:HB3	25:BC:202:ARG:HB3	1.75	0.68
26:BD:62:LYS:CD	26:BD:62:LYS:H	2.05	0.68
35:BN:115:LEU:O	35:BN:116:VAL:HG12	1.92	0.68
37:BP:21:PRO:HB2	37:BP:49:ILE:HD13	1.76	0.68
3:CD:157:ALA:HA	3:CD:160:LEU:HD12	1.76	0.68
4:CE:35:LEU:HD21	4:CE:136:VAL:HG21	1.75	0.68
18:CS:27:LYS:HB3	18:CS:27:LYS:HZ3	1.58	0.68
48:D1:26:LYS:HE2	48:D1:28:THR:HB	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1098:A:H3'	52:DI:3:LYS:HB3	1.76	0.68
23:DB:2086:U:H2'	23:DB:2087:G:C8	2.29	0.68
23:DB:2198:A:H4'	23:DB:2199:A:OP1	1.92	0.68
25:DC:63:ILE:HG23	25:DC:102:TYR:CD1	2.28	0.68
23:DB:1845:G:H4'	25:DC:268:ARG:HD3	1.76	0.68
27:DE:138:LEU:O	27:DE:143:LEU:HD21	1.94	0.68
31:DJ:68:LYS:HD2	31:DJ:72:LYS:HB2	1.76	0.68
37:DP:50:ARG:NH1	37:DP:62:LYS:HB2	2.09	0.68
2:AC:71:ARG:HH21	2:AC:74:ILE:HB	1.57	0.68
18:AS:7:GLY:H	18:AS:8:PRO:HD3	1.59	0.68
50:B3:23:HIS:CD2	50:B3:23:HIS:H	2.10	0.68
23:BB:2621:G:O2'	26:BD:164:GLN:HB2	1.94	0.68
23:BB:417:C:H2'	23:BB:418:C:C6	2.29	0.68
23:BB:2575:C:H4'	26:BD:149:ASN:N	2.09	0.68
26:BD:148:GLN:HG3	26:BD:151:THR:OG1	1.93	0.68
36:BO:48:LEU:HD23	36:BO:49:VAL:H	1.59	0.68
38:BQ:47:ARG:NH2	38:BQ:50:ARG:HB3	2.09	0.68
1:CA:695:A:H5'	10:CK:52:ARG:HH22	1.58	0.68
1:CA:947:G:H2'	1:CA:948:C:C6	2.29	0.68
20:CB:40:ILE:HG23	20:CB:200:PRO:HB2	1.76	0.68
9:CJ:9:ARG:HB2	9:CJ:99:GLN:HG3	1.75	0.68
23:DB:117:G:H5'	23:DB:126:A:H8	1.58	0.68
23:DB:162:U:H4'	23:DB:163:C:OP1	1.92	0.68
23:DB:2395:C:H2'	23:DB:2396:G:O4'	1.94	0.68
23:DB:2849:U:H4'	23:DB:2850:A:H5'	1.74	0.68
23:DB:594:U:H2'	23:DB:595:C:C6	2.28	0.68
23:DB:743:A:O2'	23:DB:744:U:H5'	1.93	0.68
25:DC:48:ILE:HG22	25:DC:49:THR:N	2.05	0.68
30:DH:54:LEU:HD23	30:DH:57:LYS:HD3	1.76	0.68
32:DK:77:ILE:HD11	32:DK:105:ARG:NH1	2.09	0.68
34:DM:16:ARG:HE	34:DM:18:ARG:NH1	1.92	0.68
44:DX:25:GLN:NE2	44:DX:29:ARG:HH21	1.92	0.68
1:AA:865:A:H2'	1:AA:866:C:C6	2.28	0.68
23:BB:1177:G:H2'	23:BB:1178:C:H6	1.59	0.68
23:BB:2198:A:H4'	23:BB:2199:A:OP1	1.94	0.68
23:BB:2395:C:H2'	23:BB:2396:G:O4'	1.94	0.68
23:BB:580:U:H2'	23:BB:581:C:C6	2.29	0.68
23:BB:2575:C:H5''	26:BD:149:ASN:HB3	1.75	0.68
52:BI:5:GLN:HG2	52:BI:6:ALA:N	2.08	0.68
1:AA:1432:G:OP1	37:BP:105:LYS:HE3	1.94	0.68
1:CA:662:U:H2'	1:CA:663:A:C8	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CE:67:ARG:HB2	4:CE:67:ARG:HH11	1.59	0.68
22:DA:61:G:H2'	22:DA:62:C:C6	2.28	0.68
23:DB:2039:U:H2'	23:DB:2040:G:C8	2.28	0.68
23:DB:2262:U:O4	43:DW:12:GLY:HA2	1.93	0.68
23:DB:2405:G:C5'	33:DL:70:LYS:HG3	2.23	0.68
23:DB:2591:C:H2'	23:DB:2592:G:H8	1.59	0.68
23:DB:2783:U:H2'	23:DB:2784:U:C6	2.29	0.68
23:DB:543:G:H2'	23:DB:544:C:H4'	1.73	0.68
27:DE:115:GLN:HB3	27:DE:117:ARG:HH11	1.58	0.68
46:DZ:59:ARG:HD3	46:DZ:63:ARG:HD3	1.76	0.68
1:AA:1252:A:H2'	1:AA:1253:G:H5''	1.75	0.68
1:AA:78:A:H2'	1:AA:79:G:H8	1.59	0.68
20:AB:116:LEU:HB3	20:AB:140:LEU:HD21	1.76	0.68
15:AP:3:THR:HG22	15:AP:66:THR:HB	1.75	0.68
23:BB:558:U:OP1	31:BJ:113:PRO:HG2	1.93	0.68
26:BD:114:LYS:HE3	26:BD:196:ALA:HB2	1.76	0.68
1:CA:1436:U:H2'	1:CA:1437:A:C8	2.28	0.68
20:CB:10:LYS:HB2	20:CB:211:LEU:HD21	1.76	0.68
4:CE:44:ARG:HG2	4:CE:72:ASN:ND2	2.09	0.68
13:CN:50:LEU:HD12	13:CN:51:PRO:HD3	1.76	0.68
47:D0:47:TYR:HB3	47:D0:52:LYS:H	1.59	0.68
23:DB:1799:G:H4'	23:DB:1800:C:O5'	1.93	0.68
23:DB:2064:C:H2'	23:DB:2065:C:H6	1.58	0.68
23:DB:794:A:H2'	23:DB:795:C:C6	2.29	0.68
27:DE:137:LYS:HZ2	27:DE:137:LYS:HA	1.57	0.68
27:DE:154:ASP:OD1	27:DE:158:PHE:HB2	1.94	0.68
29:DG:66:THR:O	29:DG:70:LEU:HD13	1.94	0.68
37:DP:26:GLU:HB3	37:DP:46:VAL:HG13	1.75	0.68
43:DW:46:ALA:HB2	43:DW:77:LYS:HD3	1.76	0.68
2:AC:57:GLU:HB2	2:AC:64:ARG:HB2	1.75	0.68
3:AD:156:ALA:O	3:AD:159:GLU:HG2	1.94	0.68
4:AE:85:LYS:HE2	4:AE:92:ARG:NH1	2.08	0.68
23:BB:240:C:H3'	23:BB:241:A:H5''	1.75	0.68
23:BB:2783:U:H2'	23:BB:2784:U:C6	2.29	0.68
23:BB:479:A:N3	23:BB:481:G:H5''	2.09	0.68
25:BC:21:PRO:O	25:BC:202:ARG:HD2	1.93	0.68
26:BD:16:THR:O	37:BP:80:VAL:HB	1.94	0.68
30:BH:89:LYS:HZ1	30:BH:123:ARG:HB3	1.59	0.68
31:BJ:13:ARG:HG3	31:BJ:13:ARG:HH11	1.59	0.68
24:BV:76:ASP:HB2	34:BM:136:MET:CG	2.24	0.68
24:BV:70:ILE:HD13	24:BV:70:ILE:N	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:865:A:H2'	1:CA:866:C:C6	2.28	0.68
5:CF:42:TRP:HB2	5:CF:59:TYR:HB2	1.76	0.68
22:DA:104:A:H2'	22:DA:105:G:O4'	1.94	0.68
23:DB:1373:A:H2'	23:DB:1374:G:O4'	1.92	0.68
23:DB:528:A:C2	23:DB:2042:A:H2'	2.29	0.68
23:DB:2471:A:HO2'	23:DB:2472:G:H8	1.40	0.68
2:AC:184:ASN:HD22	2:AC:185:THR:H	1.42	0.67
3:AD:94:GLU:HG3	3:AD:103:ARG:NH1	2.09	0.67
11:AL:35:ARG:HG3	11:AL:36:VAL:H	1.58	0.67
48:B1:7:LYS:HD3	48:B1:24:LYS:HB2	1.76	0.67
23:BB:2318:G:O2'	23:BB:2319:G:C2	2.47	0.67
23:BB:2297:A:N6	23:BB:2319:G:H5''	2.09	0.67
41:BT:30:ILE:O	41:BT:84:TYR:HA	1.94	0.67
24:BV:66:ASP:HB3	24:BV:68:LYS:HG2	1.76	0.67
4:CE:22:LYS:HB3	4:CE:29:ILE:HB	1.75	0.67
6:CG:25:PHE:HD1	6:CG:100:MET:HB2	1.59	0.67
18:CS:62:THR:HG23	18:CS:63:ASP:H	1.59	0.67
23:DB:2089:C:H2'	23:DB:2090:A:O4'	1.94	0.67
23:DB:2334:U:H3	43:DW:74:LYS:CE	2.06	0.67
23:DB:2743:U:H2'	23:DB:2744:G:H5''	1.75	0.67
37:DP:64:SER:HA	37:DP:71:ARG:CB	2.23	0.67
46:DZ:20:ASN:O	46:DZ:21:VAL:HB	1.94	0.67
8:AI:112:ARG:HB2	8:AI:112:ARG:HH11	1.57	0.67
14:AO:31:LEU:HD12	14:AO:58:MET:HB2	1.76	0.67
19:AT:67:HIS:O	19:AT:70:LYS:HG2	1.93	0.67
48:B1:29:LYS:HE3	48:B1:30:PRO:HD2	1.75	0.67
51:B4:14:CYS:HA	51:B4:26:ILE:O	1.93	0.67
23:BB:1531:C:H2'	23:BB:1532:A:C8	2.29	0.67
23:BB:1917:U:H2'	23:BB:1918:A:H5'	1.76	0.67
23:BB:598:U:H2'	23:BB:599:A:H8	1.58	0.67
25:BC:144:GLU:HB2	25:BC:146:LYS:O	1.93	0.67
25:BC:42:ARG:CG	25:BC:43:ASN:H	2.01	0.67
26:BD:98:VAL:HA	26:BD:101:PHE:CZ	2.28	0.67
26:BD:128:ARG:HH22	26:BD:130:GLN:HG3	1.60	0.67
29:BG:85:LYS:O	29:BG:86:LEU:HD22	1.94	0.67
23:BB:1163:G:H21	39:BR:91:GLN:HG3	1.58	0.67
1:CA:1026:G:H2'	1:CA:1027:C:H6	1.60	0.67
20:CB:185:ILE:HA	20:CB:199:ILE:HG13	1.76	0.67
19:CT:34:VAL:HG11	19:CT:78:LEU:HD13	1.75	0.67
47:D0:52:LYS:O	47:D0:53:VAL:HG12	1.93	0.67
23:DB:2143:C:H3'	23:DB:2144:G:H8	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:596:U:H2'	23:DB:597:G:H8	1.57	0.67
26:DD:202:ILE:HG23	26:DD:204:LYS:HE3	1.76	0.67
28:DF:151:LEU:HG	28:DF:153:ILE:HG13	1.74	0.67
34:DM:100:LYS:O	34:DM:101:VAL:HG23	1.94	0.67
43:DW:17:ALA:O	43:DW:18:LYS:HG2	1.95	0.67
1:AA:1029:U:H5''	1:AA:1030:U:H5	1.59	0.67
1:AA:1300:G:H1'	1:AA:1301:U:H5	1.59	0.67
1:AA:541:G:O2'	3:AD:39:GLN:HB2	1.93	0.67
6:AG:66:GLU:HA	6:AG:69:ARG:HE	1.59	0.67
13:AN:60:ARG:HD3	13:AN:60:ARG:N	2.05	0.67
19:AT:15:LYS:HD3	19:AT:18:LYS:HE3	1.76	0.67
23:BB:811:U:O2	23:BB:1250:G:H3'	1.95	0.67
23:BB:559:G:H1'	38:BQ:55:GLN:NE2	2.10	0.67
26:BD:89:GLU:HG2	26:BD:89:GLU:O	1.93	0.67
27:BE:128:ALA:HB3	27:BE:129:PRO:CD	2.20	0.67
38:BQ:39:ILE:O	38:BQ:43:GLN:HG3	1.94	0.67
40:BS:34:ASP:H	40:BS:37:THR:CG2	2.06	0.67
41:BT:54:GLU:N	41:BT:91:GLN:HE22	1.88	0.67
1:CA:1306:A:N6	1:CA:1331:G:O2'	2.27	0.67
2:CC:127:VAL:HG23	2:CC:128:MET:H	1.57	0.67
23:DB:1060:U:O2	23:DB:1088:A:N7	2.27	0.67
23:DB:1190:G:OP1	33:DL:38:GLN:HB2	1.94	0.67
23:DB:1201:U:H2'	23:DB:1202:G:H8	1.59	0.67
23:DB:2795:C:H2'	23:DB:2796:U:O4'	1.94	0.67
31:DJ:35:ARG:HH12	31:DJ:40:HIS:N	1.82	0.67
31:DJ:96:ARG:HD2	31:DJ:99:ARG:HH21	1.58	0.67
33:DL:18:ARG:HH12	33:DL:21:ARG:HD3	1.59	0.67
36:DO:66:GLY:N	36:DO:70:ALA:HB2	2.07	0.67
37:DP:112:ARG:H	37:DP:112:ARG:HE	1.41	0.67
37:DP:29:VAL:HG23	37:DP:47:ILE:HD11	1.76	0.67
1:AA:987:G:H2'	1:AA:988:G:H8	1.60	0.67
20:AB:131:LYS:HA	20:AB:134:LEU:HD12	1.74	0.67
8:AI:34:LEU:HD11	8:AI:47:VAL:HG21	1.76	0.67
49:B2:34:ARG:HE	49:B2:42:LEU:HD13	1.59	0.67
51:B4:11:CYS:HB3	51:B4:14:CYS:SG	2.35	0.67
23:BB:215:G:C4'	23:BB:216:A:H4'	2.25	0.67
23:BB:417:C:H2'	23:BB:418:C:H6	1.59	0.67
23:BB:596:U:H2'	23:BB:597:G:H8	1.58	0.67
52:BI:89:SER:HA	52:BI:97:VAL:CG2	2.25	0.67
23:BB:2393:U:H5''	33:BL:62:PRO:HA	1.76	0.67
35:BN:100:CYS:HB3	35:BN:111:ALA:H	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BW:35:ILE:HG12	43:BW:70:VAL:HG21	1.75	0.67
43:BW:28:GLU:N	43:BW:61:LYS:HB2	2.09	0.67
46:BZ:5:ILE:HG23	46:BZ:63:ARG:HE	1.60	0.67
1:CA:194:C:O2'	1:CA:195:A:H5'	1.94	0.67
1:CA:555:U:H2'	1:CA:556:C:H6	1.60	0.67
3:CD:87:GLU:HG2	3:CD:187:ARG:HD3	1.76	0.67
19:CT:66:ILE:HG13	19:CT:70:LYS:HE2	1.75	0.67
48:D1:46:VAL:HG22	48:D1:47:ILE:H	1.59	0.67
51:D4:26:ILE:CD1	51:D4:34:LYS:HA	2.23	0.67
23:DB:1533:C:O2'	23:DB:1534:U:H5'	1.95	0.67
23:DB:863:A:H2'	23:DB:864:G:H8	1.58	0.67
25:DC:87:SER:O	25:DC:157:ALA:HB2	1.95	0.67
25:DC:260:LYS:HG2	25:DC:262:THR:OG1	1.95	0.67
31:DJ:98:GLU:HG3	31:DJ:126:ALA:HB2	1.75	0.67
41:DT:66:LYS:H	41:DT:76:ARG:NH2	1.90	0.67
1:AA:1320:C:N4	18:AS:35:ARG:HD3	2.10	0.67
1:AA:1479:C:H2'	1:AA:1480:A:H8	1.59	0.67
9:AJ:80:THR:HG22	9:AJ:81:GLU:H	1.60	0.67
11:AL:19:ASN:O	11:AL:20:VAL:HG23	1.95	0.67
23:BB:2750:A:H3'	29:BG:3:VAL:CG1	2.24	0.67
23:BB:478:A:N1	23:BB:500:G:H4'	2.09	0.67
23:BB:794:A:H2'	23:BB:795:C:C6	2.29	0.67
30:BH:6:LEU:HA	30:BH:15:LEU:CA	2.24	0.67
36:BO:67:ASN:HA	36:BO:71:ALA:HB3	1.75	0.67
38:BQ:52:ARG:CZ	38:BQ:55:GLN:HG2	2.24	0.67
51:D4:26:ILE:CG1	51:D4:35:GLN:H	2.02	0.67
23:DB:1437:C:H2'	23:DB:1438:U:H6	1.60	0.67
23:DB:594:U:H2'	23:DB:595:C:H6	1.60	0.67
29:DG:36:LEU:HB2	29:DG:40:VAL:HG21	1.76	0.67
32:DK:43:ILE:H	32:DK:43:ILE:HD12	1.58	0.67
33:DL:58:TYR:HB3	50:D3:13:PHE:CZ	2.30	0.67
23:DB:1341:G:H5'	41:DT:61:LEU:HD21	1.76	0.67
44:DX:12:GLU:HG2	44:DX:12:GLU:O	1.94	0.67
51:B4:20:ASP:OD2	51:B4:22:VAL:HG13	1.93	0.67
23:BB:2804:U:H2'	23:BB:2805:C:H6	1.58	0.67
23:BB:2834:G:H1'	23:BB:2883:A:H61	1.58	0.67
23:BB:445:C:OP1	38:BQ:1:ALA:HB3	1.94	0.67
26:BD:138:LEU:CD2	26:BD:138:LEU:H	2.07	0.67
27:BE:164:LEU:O	27:BE:165:HIS:HB2	1.93	0.67
27:BE:60:TRP:NE1	27:BE:73:ILE:HD11	2.08	0.67
28:BF:59:ILE:HD11	28:BF:137:PHE:CD2	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BX:55:THR:HG22	44:BX:59:GLU:OE1	1.95	0.67
20:CB:206:ILE:HG22	20:CB:207:ARG:HH21	1.57	0.67
23:DB:2088:A:H2'	23:DB:2089:C:C6	2.28	0.67
23:DB:2266:A:H4'	23:DB:2267:A:C2	2.30	0.67
23:DB:479:A:O2'	23:DB:481:G:H5'	1.95	0.67
23:DB:833:A:H2'	23:DB:834:G:C8	2.30	0.67
25:DC:26:GLY:C	25:DC:27:LYS:HD2	2.15	0.67
29:DG:16:VAL:HG22	29:DG:25:ILE:HG13	1.77	0.67
34:DM:118:LYS:HB2	34:DM:118:LYS:NZ	2.09	0.67
34:DM:71:LYS:HB2	34:DM:72:PRO:HD3	1.75	0.67
23:DB:2880:C:H1'	35:DN:92:GLY:O	1.95	0.67
38:DQ:26:ALA:HA	38:DQ:30:VAL:HG23	1.76	0.67
41:DT:55:VAL:HG23	41:DT:87:LEU:H	1.59	0.67
4:AE:110:MET:HB3	4:AE:139:THR:HG21	1.76	0.67
4:AE:83:PRO:HB3	4:AE:96:GLN:HG3	1.76	0.67
15:AP:9:HIS:HE1	15:AP:29:ASN:HB2	1.59	0.67
23:BB:1250:G:O2'	23:BB:1251:C:OP1	2.12	0.67
23:BB:1486:U:H2'	23:BB:1487:U:H6	1.59	0.67
23:BB:2769:U:H2'	23:BB:2770:G:H8	1.58	0.67
23:BB:2836:U:H2'	23:BB:2837:A:H8	1.60	0.67
23:BB:282:A:H2'	23:BB:283:G:C8	2.30	0.67
23:BB:587:C:OP2	33:BL:29:LYS:HA	1.94	0.67
25:BC:79:ARG:CZ	25:BC:110:LYS:HD3	2.24	0.67
25:BC:21:PRO:HG2	25:BC:202:ARG:NH1	2.09	0.67
27:BE:17:THR:HA	27:BE:199:MET:HG3	1.76	0.67
28:BF:9:ASP:O	28:BF:10:GLU:HB2	1.93	0.67
31:BJ:39:LYS:HZ2	38:BQ:69:ARG:HD2	1.58	0.67
39:BR:46:GLU:O	39:BR:47:VAL:HB	1.95	0.67
41:BT:38:ALA:O	41:BT:42:GLU:HB2	1.95	0.67
42:BU:4:ILE:O	42:BU:27:VAL:HG11	1.95	0.67
1:CA:1296:C:H4'	1:CA:1302:C:H41	1.60	0.67
20:CB:16:GLY:HA2	20:CB:40:ILE:HG13	1.77	0.67
20:CB:27:LYS:HB3	20:CB:28:PRO:HD3	1.77	0.67
5:CF:3:HIS:HB3	5:CF:92:THR:CA	2.22	0.67
11:CL:93:ARG:H	11:CL:93:ARG:HD2	1.59	0.67
23:DB:2278:A:H62	43:DW:10:ARG:HB2	1.59	0.67
23:DB:2292:U:H2'	23:DB:2293:G:H8	1.60	0.67
23:DB:2391:G:OP2	50:D3:32:LEU:HG	1.95	0.67
23:DB:2769:U:H2'	23:DB:2770:G:H8	1.58	0.67
23:DB:598:U:H2'	23:DB:599:A:H8	1.58	0.67
25:DC:20:ASN:HB2	25:DC:202:ARG:NH1	2.10	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:63:ILE:HD12	25:DC:83:ASP:CG	2.15	0.67
26:DD:122:VAL:HA	26:DD:128:ARG:CG	2.24	0.67
41:DT:55:VAL:HG22	41:DT:56:GLU:N	2.08	0.67
1:AA:1170:A:H2'	1:AA:1171:A:O4'	1.94	0.67
1:AA:1313:U:H5''	18:AS:5:LYS:CG	2.24	0.67
20:AB:212:TYR:O	20:AB:216:VAL:HG23	1.94	0.67
5:AF:12:PRO:HG3	5:AF:54:LEU:HG	1.76	0.67
19:AT:48:LYS:O	19:AT:52:GLU:HB2	1.95	0.67
23:BB:329:G:O6	42:BU:16:LYS:HB2	1.95	0.67
23:BB:589:U:H2'	23:BB:590:A:C8	2.30	0.67
23:BB:643:A:N6	23:BB:2370:G:H1'	2.10	0.67
23:BB:751:A:H5'	40:BS:90:LYS:HE3	1.75	0.67
25:BC:238:ASN:HB3	25:BC:242:HIS:NE2	2.09	0.67
25:BC:9:SER:HB2	25:BC:202:ARG:CZ	2.25	0.67
26:BD:146:ILE:HD12	26:BD:146:ILE:N	2.10	0.67
27:BE:170:ARG:HH21	27:BE:175:ILE:HD12	1.59	0.67
29:BG:176:LYS:C	29:BG:176:LYS:HE2	2.15	0.67
32:BK:26:GLY:HA3	32:BK:30:ARG:NE	2.06	0.67
1:CA:398:U:H2'	1:CA:399:G:H8	1.60	0.67
20:CB:202:ASN:HD22	20:CB:203:ASP:H	1.43	0.67
2:CC:8:GLY:HA3	13:CN:88:MET:SD	2.35	0.67
1:CA:554:A:H5'	11:CL:25:ALA:HB1	1.77	0.67
12:CM:79:LEU:HD22	12:CM:86:ARG:HE	1.60	0.67
23:DB:1531:C:H2'	23:DB:1532:A:C8	2.29	0.67
23:DB:2900:A:H2'	23:DB:2901:C:C6	2.30	0.67
23:DB:836:G:H2'	23:DB:837:C:H6	1.58	0.67
25:DC:22:GLU:HB2	25:DC:202:ARG:CD	2.25	0.67
27:DE:6:LYS:HD3	27:DE:7:ASP:N	2.10	0.67
28:DF:39:VAL:HG22	28:DF:84:ILE:HD13	1.77	0.67
29:DG:40:VAL:HG22	29:DG:51:PHE:CE2	2.29	0.67
30:DH:6:LEU:CD1	30:DH:36:ALA:H	2.03	0.67
52:DI:27:LEU:CD2	52:DI:27:LEU:H	2.07	0.67
34:DM:80:VAL:HG12	34:DM:81:ARG:H	1.60	0.67
40:DS:73:LYS:HD3	40:DS:74:ILE:H	1.58	0.67
44:DX:51:ALA:O	44:DX:53:VAL:N	2.28	0.67
1:AA:1219:A:H2'	1:AA:1220:G:H8	1.59	0.67
1:AA:203:G:H1'	1:AA:465:A:H62	1.59	0.67
1:AA:764:C:H3'	1:AA:765:G:N2	2.06	0.67
6:AG:149:ALA:HB1	10:AK:58:THR:HG21	1.77	0.67
8:AI:82:ILE:O	8:AI:86:LEU:HD22	1.95	0.67
23:BB:1210:G:H5'	23:BB:1212:G:O4'	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2039:U:H2'	23:BB:2040:G:C8	2.28	0.67
23:BB:2294:G:O2'	23:BB:2295:C:H5'	1.95	0.67
23:BB:721:A:H2'	23:BB:722:A:C8	2.29	0.67
52:BI:77:VAL:HA	52:BI:80:LYS:HE2	1.76	0.67
23:BB:1007:C:H5''	31:BJ:37:ARG:HH21	1.59	0.67
34:BM:35:ALA:O	34:BM:36:VAL:HG22	1.95	0.67
46:BZ:36:VAL:CA	46:BZ:42:PRO:HB3	2.23	0.67
1:CA:763:G:H2'	1:CA:764:C:H6	1.60	0.67
12:CM:39:ALA:O	12:CM:42:VAL:HG22	1.94	0.67
23:DB:1568:G:OP1	25:DC:61:TYR:HB2	1.95	0.67
26:DD:130:GLN:HB3	26:DD:140:HIS:HA	1.77	0.67
26:DD:48:ILE:HA	26:DD:80:TRP:HB3	1.77	0.67
33:DL:34:GLY:HA3	39:DR:85:LYS:HG2	1.77	0.67
33:DL:39:LYS:H	33:DL:39:LYS:HZ3	1.42	0.67
37:DP:7:LEU:HD23	37:DP:11:GLN:HE21	1.60	0.67
39:DR:47:VAL:CG1	39:DR:49:ILE:HG12	2.25	0.67
43:DW:42:THR:HB	43:DW:75:ASN:CB	2.24	0.67
1:AA:1218:C:H2'	1:AA:1219:A:H8	1.59	0.67
1:AA:586:C:O2'	1:AA:587:G:H5'	1.94	0.67
5:AF:53:LYS:HA	5:AF:53:LYS:NZ	2.09	0.67
7:AH:29:SER:HB3	7:AH:32:LYS:HG3	1.77	0.67
11:AL:20:VAL:HG13	11:AL:94:TYR:OH	1.95	0.67
16:AQ:29:LYS:HD3	16:AQ:35:LYS:N	2.10	0.67
10:AK:110:THR:HG21	21:AU:4:LYS:HD2	1.77	0.67
50:B3:41:ARG:HB3	50:B3:43:LEU:HD22	1.77	0.67
23:BB:1410:G:H2'	23:BB:1411:U:C6	2.29	0.67
23:BB:2537:U:H2'	23:BB:2538:C:H6	1.60	0.67
23:BB:442:G:N2	27:BE:46:GLN:NE2	2.44	0.67
34:BM:3:GLN:O	34:BM:5:LYS:N	2.28	0.67
35:BN:114:GLU:HG2	35:BN:115:LEU:H	1.59	0.67
1:CA:555:U:H2'	1:CA:556:C:C6	2.30	0.67
1:CA:673:A:H2'	1:CA:674:G:C8	2.28	0.67
1:CA:918:A:H2'	1:CA:919:A:C8	2.30	0.67
20:CB:103:TRP:O	20:CB:107:ARG:HG3	1.95	0.67
9:CJ:12:ALA:HB3	9:CJ:18:ILE:HB	1.75	0.67
1:CA:1202:U:H1'	13:CN:68:ARG:HD2	1.77	0.67
22:DA:26:C:H2'	22:DA:27:C:C6	2.29	0.67
25:DC:107:LYS:CE	25:DC:108:GLY:H	2.05	0.67
27:DE:138:LEU:HD22	27:DE:187:VAL:HG11	1.75	0.67
39:DR:74:ILE:O	39:DR:75:VAL:HG13	1.94	0.67
1:AA:1065:U:H1'	1:AA:1066:C:OP2	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1193:G:O2'	1:AA:1194:U:H5'	1.95	0.66
1:AA:398:U:H2'	1:AA:399:G:H8	1.60	0.66
1:AA:555:U:H2'	1:AA:556:C:C6	2.30	0.66
20:AB:96:LEU:HB2	20:AB:99:MET:HE3	1.76	0.66
9:AJ:10:LEU:HG	9:AJ:98:VAL:HG12	1.76	0.66
13:AN:50:LEU:HD23	13:AN:51:PRO:HD3	1.77	0.66
23:BB:833:A:H2'	23:BB:834:G:C8	2.30	0.66
26:BD:34:VAL:HG13	26:BD:89:GLU:OE2	1.94	0.66
35:BN:72:ASP:HB3	35:BN:74:GLU:HG3	1.76	0.66
1:CA:678:U:H2'	1:CA:679:C:C6	2.30	0.66
1:CA:678:U:H2'	1:CA:679:C:H6	1.59	0.66
1:CA:408:A:OP1	3:CD:111:ALA:HB3	1.95	0.66
4:CE:87:VAL:HG22	4:CE:88:HIS:N	2.09	0.66
5:CF:68:GLN:HA	5:CF:71:ILE:HG22	1.77	0.66
6:CG:70:PRO:HA	6:CG:141:HIS:HE1	1.59	0.66
1:CA:538:G:OP2	11:CL:111:GLN:HB2	1.94	0.66
11:CL:65:TYR:HB2	11:CL:92:VAL:HG11	1.77	0.66
23:DB:2723:C:H5'	35:DN:4:ARG:CZ	2.25	0.66
23:DB:2836:U:H2'	23:DB:2837:A:H8	1.60	0.66
23:DB:1799:G:H5''	25:DC:255:LYS:HE3	1.76	0.66
26:DD:122:VAL:HG22	26:DD:128:ARG:HG3	1.77	0.66
31:DJ:37:ARG:HH21	31:DJ:46:PRO:HB3	1.59	0.66
42:DU:43:LYS:HB2	42:DU:43:LYS:HZ2	1.60	0.66
1:AA:662:U:H2'	1:AA:663:A:C8	2.30	0.66
1:AA:763:G:H2'	1:AA:764:C:H6	1.59	0.66
16:AQ:75:VAL:HG23	16:AQ:76:ARG:H	1.59	0.66
23:BB:2527:C:H1'	51:B4:1:MET:HB3	1.77	0.66
23:BB:1082:U:N3	23:BB:1086:A:C6	2.63	0.66
23:BB:2147:A:H5'	23:BB:2148:G:C4'	2.26	0.66
23:BB:2591:C:H2'	23:BB:2592:G:H8	1.59	0.66
35:BN:97:ILE:CD1	35:BN:113:ILE:HD13	2.24	0.66
39:BR:5:PHE:HA	39:BR:40:MET:CG	2.25	0.66
46:BZ:55:GLY:HA2	46:BZ:63:ARG:HG3	1.78	0.66
1:CA:909:A:H2'	1:CA:910:C:O4'	1.95	0.66
8:CI:98:ARG:HA	8:CI:103:VAL:HG22	1.75	0.66
9:CJ:48:ARG:HB3	9:CJ:66:GLU:HG3	1.78	0.66
9:CJ:8:ILE:HA	9:CJ:100:ILE:HG22	1.76	0.66
11:CL:54:VAL:HG11	11:CL:79:ILE:HD11	1.77	0.66
15:CP:61:VAL:HA	15:CP:65:ALA:HB3	1.77	0.66
21:CU:48:LYS:HA	21:CU:51:ALA:HB3	1.77	0.66
48:D1:16:THR:HG22	48:D1:47:ILE:HD12	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:D4:34:LYS:HE2	51:D4:36:ARG:NH2	2.10	0.66
23:DB:1410:G:H2'	23:DB:1411:U:C6	2.29	0.66
23:DB:1486:U:H2'	23:DB:1487:U:H6	1.59	0.66
23:DB:532:A:H4'	23:DB:533:G:C8	2.30	0.66
30:DH:87:GLU:HB2	30:DH:89:LYS:NZ	2.08	0.66
31:DJ:132:HIS:HB3	31:DJ:136:GLN:OE1	1.95	0.66
32:DK:19:VAL:HG12	32:DK:43:ILE:HA	1.77	0.66
37:DP:36:LYS:HG2	37:DP:37:LYS:HE3	1.76	0.66
37:DP:55:HIS:C	37:DP:57:ALA:H	1.96	0.66
42:DU:17:ASP:OD2	42:DU:20:LYS:HB2	1.95	0.66
42:DU:82:VAL:HB	42:DU:94:PHE:HB3	1.76	0.66
1:AA:1116:U:O2'	1:AA:1117:A:H5'	1.95	0.66
1:AA:678:U:H2'	1:AA:679:C:C6	2.30	0.66
23:BB:1212:G:O2'	23:BB:1236:G:N2	2.27	0.66
23:BB:1437:C:H2'	23:BB:1438:U:H6	1.59	0.66
23:BB:2471:A:HO2'	23:BB:2472:G:H8	1.37	0.66
23:BB:2511:U:O2'	26:BD:144:GLY:HA3	1.94	0.66
27:BE:117:ARG:H	27:BE:118:LEU:HD23	1.60	0.66
31:BJ:56:VAL:HB	31:BJ:124:VAL:HA	1.77	0.66
35:BN:34:ILE:HD13	35:BN:34:ILE:N	2.10	0.66
40:BS:103:ILE:HG13	40:BS:105:VAL:HG23	1.76	0.66
1:CA:1078:U:H4'	4:CE:137:ARG:NH1	2.10	0.66
1:CA:1270:G:H2'	1:CA:1271:A:H8	1.61	0.66
1:CA:731:G:H5'	1:CA:766:A:H4'	1.78	0.66
6:CG:2:ARG:HB3	6:CG:2:ARG:NH1	2.10	0.66
8:CI:29:ILE:HA	8:CI:64:ILE:HB	1.76	0.66
23:DB:2286:G:H1	48:D1:23:THR:HG21	1.60	0.66
23:DB:1456:G:O2'	23:DB:1457:U:H5'	1.95	0.66
23:DB:2243:U:H2'	23:DB:2244:U:C6	2.30	0.66
26:DD:125:TRP:HB2	26:DD:160:LYS:CD	2.25	0.66
39:DR:6:GLN:HG2	39:DR:7:SER:H	1.60	0.66
42:DU:42:LYS:H	42:DU:57:ILE:HD12	1.58	0.66
43:DW:31:LEU:HD22	43:DW:39:GLN:HE22	1.61	0.66
1:AA:1008:U:H2'	1:AA:1009:U:C4'	2.25	0.66
1:AA:1163:A:H2'	1:AA:1164:G:H8	1.60	0.66
1:AA:72:A:H2'	1:AA:73:C:H6	1.60	0.66
3:AD:2:ARG:HG3	3:AD:114:ARG:CZ	2.26	0.66
2:AC:5:HIS:CD2	13:AN:88:MET:HB3	2.30	0.66
23:BB:161:A:H62	23:BB:165:A:H61	1.41	0.66
23:BB:1714:U:H3'	23:BB:1715:G:H5'	1.77	0.66
27:BE:148:ILE:HB	27:BE:185:LYS:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:BG:40:VAL:HG21	29:BG:53:PRO:HA	1.76	0.66
31:BJ:23:LYS:HD2	31:BJ:23:LYS:H	1.61	0.66
43:BW:66:VAL:HG22	43:BW:67:LYS:N	2.11	0.66
1:CA:16:A:O2'	1:CA:17:U:H5'	1.96	0.66
1:CA:376:G:H5''	15:CP:5:ARG:HB2	1.76	0.66
9:CJ:37:ARG:HB2	9:CJ:76:ILE:HA	1.75	0.66
23:DB:1082:U:N3	23:DB:1086:A:C6	2.62	0.66
23:DB:2323:G:O2'	23:DB:2324:U:H5'	1.94	0.66
23:DB:2531:A:OP1	29:DG:176:LYS:HA	1.95	0.66
23:DB:857:G:H2'	23:DB:858:G:H5'	1.78	0.66
25:DC:74:PRO:HB2	25:DC:96:LYS:HG3	1.78	0.66
26:DD:4:LEU:HD22	26:DD:51:THR:HB	1.76	0.66
27:DE:105:LEU:O	27:DE:108:ILE:HG23	1.95	0.66
29:DG:8:VAL:HB	29:DG:49:LEU:HB2	1.77	0.66
40:DS:23:LEU:C	40:DS:24:ILE:HD13	2.15	0.66
41:DT:72:GLN:C	41:DT:73:ARG:HD2	2.16	0.66
24:DV:4:ILE:O	24:DV:63:ILE:HA	1.95	0.66
2:AC:77:GLY:HA3	2:AC:82:ASP:H	1.61	0.66
18:AS:49:ALA:HA	18:AS:57:VAL:O	1.96	0.66
23:BB:1381:G:C2'	23:BB:1382:G:H5'	2.26	0.66
23:BB:1843:C:O4'	25:BC:250:GLN:HG2	1.95	0.66
23:BB:469:G:OP2	27:BE:55:SER:HB2	1.96	0.66
27:BE:5:LEU:H	27:BE:5:LEU:HD12	1.60	0.66
52:BI:11:GLN:HA	52:BI:55:PRO:HA	1.76	0.66
33:BL:118:THR:H	33:BL:119:PRO:CD	2.06	0.66
34:BM:8:LYS:HZ2	34:BM:70:ASP:N	1.92	0.66
37:BP:29:VAL:O	37:BP:31:VAL:HG23	1.95	0.66
23:BB:2270:A:H4'	43:BW:18:LYS:HG2	1.78	0.66
1:CA:239:U:H6	1:CA:239:U:C5'	2.08	0.66
1:CA:56:U:H2'	1:CA:57:G:H8	1.61	0.66
1:CA:812:G:HO2'	1:CA:813:U:H6	1.42	0.66
6:CG:113:LYS:HE2	6:CG:113:LYS:HA	1.75	0.66
49:D2:13:ASN:OD1	49:D2:17:GLY:HA3	1.96	0.66
23:DB:1098:A:H3'	52:DI:3:LYS:C	2.15	0.66
23:DB:2147:A:H2'	23:DB:2147:A:N3	2.11	0.66
23:DB:2758:A:C2'	23:DB:2759:G:H5'	2.25	0.66
26:DD:114:LYS:HB2	26:DD:114:LYS:NZ	2.10	0.66
26:DD:35:THR:HB	26:DD:48:ILE:HB	1.77	0.66
29:DG:153:PRO:HA	29:DG:159:LYS:O	1.94	0.66
29:DG:4:ALA:O	29:DG:5:LYS:HB2	1.93	0.66
34:DM:5:LYS:HD3	34:DM:71:LYS:HG2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:90:ALA:HB3	37:DP:112:ARG:N	2.11	0.66
39:DR:6:GLN:HG2	39:DR:7:SER:N	2.10	0.66
24:DV:9:ARG:NH1	24:DV:12:GLN:HA	2.09	0.66
43:DW:35:ILE:HB	43:DW:67:LYS:HZ2	1.60	0.66
1:AA:1391:U:H2'	1:AA:1392:G:C8	2.31	0.66
1:AA:678:U:H2'	1:AA:679:C:H6	1.59	0.66
1:AA:922:G:H2'	1:AA:923:A:C8	2.31	0.66
15:AP:46:LYS:H	15:AP:46:LYS:CD	2.08	0.66
23:BB:163:C:H2'	23:BB:164:C:O4'	1.95	0.66
23:BB:320:A:H4'	23:BB:322:A:N7	2.11	0.66
23:BB:448:U:H1'	27:BE:79:ARG:HG3	1.76	0.66
23:BB:545:U:H2'	23:BB:547:A:OP2	1.94	0.66
35:BN:33:ILE:HD13	35:BN:33:ILE:N	2.10	0.66
37:BP:99:LEU:HD12	37:BP:101:GLU:OE2	1.96	0.66
42:BU:5:ARG:HG2	42:BU:6:ARG:H	1.60	0.66
43:BW:46:ALA:H	43:BW:75:ASN:CG	1.99	0.66
1:CA:1053:G:HO2'	1:CA:1199:U:H5	1.42	0.66
1:CA:1461:G:H2'	1:CA:1462:C:H6	1.59	0.66
1:CA:205:A:H2'	1:CA:206:C:H6	1.61	0.66
2:CC:176:THR:OG1	2:CC:179:ALA:HB2	1.96	0.66
23:DB:1099:G:C5'	52:DI:3:LYS:N	2.59	0.66
23:DB:580:U:H2'	23:DB:581:C:C6	2.29	0.66
23:DB:928:A:H1'	45:DY:1:ALA:HA	1.77	0.66
23:DB:943:A:P	33:DL:40:SER:HA	2.35	0.66
25:DC:75:ALA:HB1	25:DC:93:VAL:HG13	1.78	0.66
26:DD:129:THR:HG23	26:DD:130:GLN:H	1.60	0.66
36:DO:25:ARG:CG	36:DO:94:ARG:HH22	2.08	0.66
42:DU:71:ILE:HD13	42:DU:72:PHE:N	2.11	0.66
45:DY:2:LYS:CB	45:DY:37:ARG:HB2	2.25	0.66
1:AA:1110:A:H2'	1:AA:1111:A:H8	1.61	0.66
1:AA:539:A:H2'	1:AA:540:G:C8	2.31	0.66
1:AA:555:U:H2'	1:AA:556:C:H6	1.60	0.66
1:AA:731:G:H5'	1:AA:766:A:H4'	1.78	0.66
3:AD:187:ARG:HA	3:AD:190:LEU:HD22	1.78	0.66
1:AA:37:U:P	11:AL:119:LYS:HB2	2.35	0.66
15:AP:51:ARG:C	15:AP:52:LEU:HD22	2.16	0.66
23:BB:1456:G:O2'	23:BB:1457:U:H5'	1.95	0.66
25:BC:257:ARG:HA	25:BC:257:ARG:NH1	2.11	0.66
34:BM:100:LYS:HE3	34:BM:100:LYS:N	2.10	0.66
40:BS:70:LYS:HD2	40:BS:110:ARG:HA	1.78	0.66
1:CA:1286:U:O2'	1:CA:1287:A:H5''	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:312:C:H2'	1:CA:313:A:C8	2.31	0.66
9:CJ:41:PRO:HA	9:CJ:72:ARG:HD3	1.77	0.66
10:CK:12:ARG:C	10:CK:13:LYS:HD2	2.15	0.66
10:CK:124:LYS:HD3	21:CU:34:ARG:HD3	1.76	0.66
23:DB:1083:U:H1'	23:DB:1086:A:H61	1.60	0.66
23:DB:1098:A:C4	52:DI:3:LYS:O	2.49	0.66
23:DB:564:C:O2'	23:DB:565:C:H5'	1.95	0.66
23:DB:589:U:H2'	23:DB:590:A:C8	2.30	0.66
25:DC:208:GLY:HA2	25:DC:212:TRP:HB3	1.78	0.66
52:DI:41:PHE:O	52:DI:45:THR:HG23	1.96	0.66
34:DM:16:ARG:HE	34:DM:18:ARG:HH12	1.42	0.66
34:DM:71:LYS:NZ	34:DM:91:TYR:HB3	2.10	0.66
43:DW:30:VAL:HG12	43:DW:31:LEU:H	1.61	0.66
43:DW:44:PHE:HB3	43:DW:77:LYS:HB3	1.76	0.66
46:DZ:30:HIS:CE1	46:DZ:49:ARG:HH12	2.14	0.66
1:AA:16:A:O2'	1:AA:17:U:H5'	1.95	0.66
1:AA:239:U:C5'	1:AA:239:U:H6	2.08	0.66
3:AD:56:GLU:HG2	3:AD:195:ASN:HD22	1.60	0.66
48:B1:15:GLY:H	48:B1:47:ILE:HD12	1.60	0.66
23:BB:1219:U:H2'	23:BB:1220:G:C8	2.30	0.66
26:BD:35:THR:HB	26:BD:48:ILE:HB	1.76	0.66
28:BF:34:THR:HG23	28:BF:89:THR:HG22	1.76	0.66
33:BL:78:ARG:HD3	33:BL:81:ASP:HB2	1.78	0.66
34:BM:22:GLN:HG2	34:BM:23:GLY:N	2.11	0.66
44:BX:49:ASP:HB3	44:BX:52:ARG:HH21	1.59	0.66
1:CA:1024:G:O2'	1:CA:1025:U:H5'	1.94	0.66
1:CA:1513:A:H2'	1:CA:1514:G:H8	1.60	0.66
1:CA:539:A:H2'	1:CA:540:G:C8	2.30	0.66
1:CA:586:C:O2'	1:CA:587:G:H5'	1.95	0.66
20:CB:137:THR:HA	20:CB:140:LEU:HD12	1.77	0.66
5:CF:47:LEU:HD21	5:CF:57:ALA:HB2	1.77	0.66
12:CM:73:SER:HB3	12:CM:77:LYS:HE2	1.77	0.66
22:DA:47:C:H5'	36:DO:97:PHE:CZ	2.31	0.66
23:DB:143:C:H6	23:DB:143:C:O5'	1.79	0.66
23:DB:235:U:H2'	23:DB:236:C:C6	2.30	0.66
23:DB:2591:C:H2'	23:DB:2592:G:C8	2.31	0.66
23:DB:289:G:H2'	23:DB:290:U:H6	1.58	0.66
23:DB:414:C:H2'	23:DB:415:A:H8	1.59	0.66
23:DB:65:U:H2'	23:DB:66:C:H6	1.61	0.66
39:DR:41:ILE:HG13	39:DR:43:ASN:HD22	1.61	0.66
40:DS:84:ARG:HH21	40:DS:98:LYS:NZ	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DW:42:THR:H	43:DW:65:LYS:HA	1.61	0.66
1:AA:1300:G:H1'	1:AA:1301:U:C5	2.31	0.66
23:BB:142:A:H2'	23:BB:143:C:C6	2.31	0.66
23:BB:2246:G:H2'	23:BB:2247:A:C8	2.30	0.66
23:BB:2461:A:H2'	23:BB:2462:C:H6	1.59	0.66
29:BG:86:LEU:HB2	29:BG:130:ILE:HB	1.76	0.66
52:BI:91:LYS:HB2	52:BI:94:LYS:HD2	1.78	0.66
31:BJ:24:THR:HB	31:BJ:27:ARG:HB2	1.78	0.66
35:BN:28:LEU:H	35:BN:28:LEU:HD12	1.61	0.66
1:CA:1462:C:H4'	37:DP:110:LYS:HE2	1.77	0.66
1:CA:997:U:H2'	1:CA:998:C:C6	2.30	0.66
6:CG:42:VAL:O	6:CG:46:LEU:HB2	1.96	0.66
10:CK:12:ARG:N	10:CK:76:TYR:HA	2.10	0.66
12:CM:78:ARG:O	12:CM:82:LEU:HG	1.95	0.66
23:DB:304:U:H2'	23:DB:305:C:C6	2.31	0.66
26:DD:95:SER:C	26:DD:96:ILE:HG13	2.16	0.66
44:DX:11:VAL:HG12	44:DX:13:GLU:H	1.60	0.66
1:AA:56:U:H2'	1:AA:57:G:H8	1.61	0.66
2:AC:154:GLY:HA3	2:AC:162:ALA:HB1	1.75	0.66
7:AH:55:LYS:HA	7:AH:55:LYS:HE3	1.77	0.66
9:AJ:6:ILE:HA	9:AJ:102:LEU:HG	1.78	0.66
11:AL:106:VAL:HG22	11:AL:116:TYR:HB3	1.76	0.66
11:AL:42:LYS:HE3	11:AL:90:PRO:HD3	1.78	0.66
23:BB:1645:G:H5''	23:BB:1646:C:H5'	1.78	0.66
23:BB:564:C:O2'	23:BB:565:C:H5'	1.95	0.66
23:BB:594:U:H2'	23:BB:595:C:H6	1.60	0.66
25:BC:204:LEU:HB3	25:BC:210:ALA:HB1	1.78	0.66
26:BD:170:VAL:HG13	26:BD:171:THR:N	2.08	0.66
28:BF:125:GLY:HA3	28:BF:159:ALA:HB3	1.77	0.66
30:BH:26:ALA:HA	30:BH:30:LEU:HB2	1.77	0.66
23:BB:1060:U:OP2	52:BI:74:PRO:HA	1.95	0.66
31:BJ:71:ASP:CG	31:BJ:72:LYS:N	2.47	0.66
33:BL:142:ILE:HG13	33:BL:143:GLU:N	2.10	0.66
37:BP:38:ARG:HG3	37:BP:39:LEU:H	1.61	0.66
39:BR:23:GLU:HA	39:BR:96:VAL:HG13	1.77	0.66
1:CA:105:G:H2'	1:CA:106:C:C6	2.31	0.66
1:CA:1132:C:H2'	1:CA:1133:G:C8	2.30	0.66
1:CA:1422:G:O2'	1:CA:1423:G:H5'	1.95	0.66
1:CA:960:U:O2'	1:CA:1223:C:H5'	1.96	0.66
4:CE:81:GLN:HE22	4:CE:148:SER:HA	1.61	0.66
5:CF:35:LYS:O	5:CF:64:VAL:HG13	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CI:14:SER:HA	8:CI:68:GLY:O	1.96	0.66
16:CQ:45:VAL:HG21	16:CQ:60:ILE:HG21	1.77	0.66
23:DB:813:U:H2'	23:DB:814:C:C6	2.31	0.66
23:DB:728:G:H1'	25:DC:16:VAL:HG21	1.77	0.66
31:DJ:97:PRO:O	31:DJ:100:VAL:HG12	1.96	0.66
37:DP:71:ARG:O	37:DP:72:VAL:HG22	1.96	0.66
42:DU:22:GLY:HA3	42:DU:35:VAL:HG23	1.78	0.66
24:DV:53:LYS:HZ2	24:DV:55:GLU:HG2	1.60	0.66
46:DZ:27:THR:O	46:DZ:28:VAL:HG13	1.96	0.66
13:AN:30:ILE:HB	13:AN:44:VAL:HG11	1.78	0.65
5:AF:86:ARG:NH1	17:AR:63:TYR:HB3	2.11	0.65
23:BB:1534:U:HO2'	23:BB:1535:A:H8	1.42	0.65
23:BB:2049:G:O2'	23:BB:2050:C:H5'	1.96	0.65
23:BB:2758:A:H2'	23:BB:2759:G:O4'	1.96	0.65
23:BB:704:G:H1'	23:BB:727:A:H61	1.60	0.65
23:BB:857:G:C2'	23:BB:858:G:H5'	2.26	0.65
23:BB:869:G:H2'	23:BB:870:U:H6	1.61	0.65
23:BB:691:C:H4'	25:BC:42:ARG:NE	2.11	0.65
26:BD:167:ASN:H	26:BD:167:ASN:HD22	1.43	0.65
38:BQ:85:ALA:CB	38:BQ:88:GLU:HB2	2.26	0.65
41:BT:77:ARG:HG2	41:BT:78:SER:H	1.60	0.65
1:CA:1119:C:H2'	1:CA:1120:C:H6	1.60	0.65
1:CA:376:G:H4'	15:CP:5:ARG:HD3	1.78	0.65
2:CC:166:TRP:CG	2:CC:167:TYR:N	2.65	0.65
4:CE:131:ASN:O	4:CE:135:VAL:HG23	1.95	0.65
13:CN:11:LYS:O	13:CN:15:LEU:HG	1.96	0.65
51:D4:2:LYS:HB2	51:D4:2:LYS:NZ	2.10	0.65
23:DB:2295:C:O2'	23:DB:2296:U:H5'	1.96	0.65
25:DC:161:VAL:HG12	25:DC:173:LEU:HD22	1.77	0.65
25:DC:20:ASN:HB2	25:DC:202:ARG:HD3	1.78	0.65
31:DJ:18:VAL:O	31:DJ:56:VAL:HA	1.96	0.65
34:DM:55:ARG:O	34:DM:57:VAL:HG23	1.96	0.65
35:DN:45:ARG:NH2	35:DN:113:ILE:HG23	2.04	0.65
35:DN:45:ARG:CZ	35:DN:95:THR:HB	2.26	0.65
24:DV:2:PHE:HB2	24:DV:61:LEU:HD22	1.78	0.65
44:DX:31:GLN:HE21	44:DX:31:GLN:HA	1.59	0.65
1:AA:105:G:H2'	1:AA:106:C:C6	2.31	0.65
1:AA:312:C:H2'	1:AA:313:A:C8	2.30	0.65
9:AJ:37:ARG:NE	9:AJ:77:VAL:HG11	2.12	0.65
21:AU:39:LYS:N	21:AU:40:PRO:HD2	2.10	0.65
23:BB:1533:C:O2'	23:BB:1534:U:H5'	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2027:G:O2'	23:BB:2028:U:H5'	1.96	0.65
23:BB:6:A:H2'	23:BB:7:G:H8	1.62	0.65
23:BB:1076:C:H4'	52:BI:94:LYS:NZ	2.11	0.65
35:BN:49:GLU:HB2	35:BN:50:PRO:HD3	1.77	0.65
45:BY:6:ILE:HD13	45:BY:6:ILE:H	1.61	0.65
1:CA:1060:U:H2'	1:CA:1061:G:H8	1.61	0.65
1:CA:1320:C:OP1	18:CS:69:LYS:HE3	1.96	0.65
1:CA:37:U:OP1	11:CL:120:ARG:HB3	1.96	0.65
1:CA:501:C:H2'	1:CA:502:A:C8	2.31	0.65
4:CE:87:VAL:HG13	4:CE:88:HIS:N	2.10	0.65
51:D4:3:VAL:HG12	51:D4:4:ARG:N	2.07	0.65
23:DB:2443:C:H2'	23:DB:2444:G:C8	2.31	0.65
26:DD:146:ILE:HG12	26:DD:155:VAL:HG13	1.76	0.65
31:DJ:41:LYS:HZ3	31:DJ:45:THR:HA	1.62	0.65
39:DR:74:ILE:HG13	39:DR:76:LYS:HG3	1.78	0.65
23:DB:2012:G:H4'	40:DS:96:ILE:HD11	1.78	0.65
43:DW:19:ARG:NH1	43:DW:19:ARG:HB2	2.11	0.65
1:AA:93:U:H3'	1:AA:94:G:H5''	1.78	0.65
23:BB:1590:A:H2'	23:BB:1591:A:C8	2.32	0.65
25:BC:139:THR:HA	25:BC:193:GLU:OE1	1.96	0.65
25:BC:21:PRO:O	25:BC:22:GLU:HB2	1.96	0.65
43:BW:44:PHE:O	43:BW:75:ASN:HB3	1.96	0.65
44:BX:9:LYS:HE3	44:BX:11:VAL:HG13	1.78	0.65
1:CA:389:A:H3'	1:CA:390:U:H6	1.61	0.65
1:CA:518:C:H2'	1:CA:530:G:C8	2.31	0.65
20:CB:208:ALA:O	20:CB:211:LEU:HB3	1.96	0.65
3:CD:30:LYS:HB2	3:CD:30:LYS:HZ2	1.61	0.65
10:CK:88:PRO:CD	21:CU:28:LEU:HD13	2.26	0.65
23:DB:1197:G:H2'	23:DB:1198:U:H6	1.62	0.65
23:DB:1381:G:C2'	23:DB:1382:G:H5'	2.26	0.65
23:DB:1590:A:H2'	23:DB:1591:A:C8	2.32	0.65
23:DB:2027:G:O2'	23:DB:2028:U:H5'	1.96	0.65
23:DB:2307:G:N3	23:DB:2307:G:H2'	2.11	0.65
23:DB:181:A:H1'	23:DB:435:C:H5'	1.78	0.65
23:DB:899:A:H3'	23:DB:900:A:C8	2.30	0.65
26:DD:146:ILE:N	26:DD:146:ILE:HD12	2.11	0.65
26:DD:151:THR:O	26:DD:153:GLY:N	2.28	0.65
26:DD:46:ARG:HA	26:DD:82:PHE:HA	1.79	0.65
34:DM:40:ARG:HA	34:DM:92:TRP:HE1	1.62	0.65
37:DP:50:ARG:HH12	37:DP:62:LYS:HB2	1.59	0.65
46:DZ:5:ILE:O	46:DZ:51:VAL:HG13	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1381:U:O2'	1:AA:1382:C:H5'	1.97	0.65
20:AB:52:ALA:O	20:AB:56:LEU:HD13	1.96	0.65
11:AL:37:TYR:HB2	11:AL:51:VAL:HG23	1.79	0.65
47:B0:2:VAL:HG22	47:B0:3:GLN:N	2.11	0.65
23:BB:1112:G:H4'	29:BG:2:ARG:HG2	1.78	0.65
23:BB:2008:C:H2'	23:BB:2009:A:C8	2.31	0.65
23:BB:710:U:H2'	23:BB:711:G:H8	1.61	0.65
23:BB:936:A:H2'	23:BB:937:C:C6	2.30	0.65
23:BB:95:A:H4'	44:BX:40:SER:HB3	1.78	0.65
25:BC:10:PRO:HD2	25:BC:202:ARG:NH1	2.11	0.65
26:BD:92:VAL:HG22	26:BD:94:GLN:HB2	1.78	0.65
27:BE:130:LYS:HA	27:BE:130:LYS:HE2	1.77	0.65
28:BF:110:ILE:HG22	28:BF:112:ASP:H	1.61	0.65
24:BV:76:ASP:HB2	34:BM:136:MET:HG3	1.78	0.65
34:BM:3:GLN:CD	34:BM:5:LYS:H	1.98	0.65
42:BU:7:ASP:CA	42:BU:24:VAL:HA	2.26	0.65
44:BX:5:GLU:H	44:BX:7:ARG:HE	1.44	0.65
1:CA:1343:G:H2'	1:CA:1344:C:C6	2.31	0.65
8:CI:94:ARG:HD3	8:CI:97:LEU:HD12	1.79	0.65
13:CN:31:SER:HA	13:CN:40:ARG:HA	1.77	0.65
47:D0:27:LEU:N	47:D0:27:LEU:HD22	2.11	0.65
47:D0:41:HIS:HB3	47:D0:47:TYR:H	1.62	0.65
33:DL:63:LYS:HB2	50:D3:26:ALA:HB2	1.78	0.65
23:DB:1714:U:H3'	23:DB:1715:G:H5'	1.77	0.65
23:DB:225:C:O2'	23:DB:226:A:H5'	1.97	0.65
23:DB:2461:A:H2'	23:DB:2462:C:H6	1.60	0.65
23:DB:4:U:H2'	23:DB:5:A:C8	2.31	0.65
25:DC:32:LEU:HA	25:DC:81:GLU:OE2	1.96	0.65
26:DD:122:VAL:HG21	26:DD:141:ARG:HD3	1.78	0.65
29:DG:39:ALA:O	29:DG:41:GLU:HG3	1.96	0.65
38:DQ:9:ALA:O	38:DQ:12:ARG:HB3	1.97	0.65
42:DU:59:GLU:HG2	42:DU:60:LYS:N	2.10	0.65
1:AA:967:C:OP1	1:AA:969:A:H5'	1.97	0.65
2:AC:88:LYS:HD3	2:AC:88:LYS:O	1.97	0.65
13:AN:12:ARG:HH11	13:AN:60:ARG:NH1	1.94	0.65
15:AP:42:ILE:HG22	15:AP:43:ALA:N	2.12	0.65
23:BB:1856:U:H2'	23:BB:1857:G:O4'	1.96	0.65
23:BB:2591:C:H2'	23:BB:2592:G:C8	2.31	0.65
37:BP:52:ARG:HG2	37:BP:54:LEU:HB2	1.77	0.65
38:BQ:47:ARG:O	38:BQ:51:GLN:HB3	1.97	0.65
40:BS:17:VAL:HA	40:BS:43:ALA:HB1	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:44:HIS:HB2	42:BU:56:GLY:N	2.11	0.65
1:CA:1132:C:H2'	1:CA:1133:G:H8	1.61	0.65
1:CA:190:A:O5'	1:CA:190:A:H8	1.80	0.65
20:CB:153:MET:SD	20:CB:157:PRO:HD3	2.35	0.65
12:CM:48:SER:H	12:CM:51:GLN:HE22	1.43	0.65
18:CS:62:THR:H	18:CS:65:MET:HE3	1.62	0.65
50:D3:49:VAL:HG22	50:D3:50:SER:N	2.08	0.65
22:DA:5:U:H2'	22:DA:6:G:H8	1.61	0.65
23:DB:1360:G:H2'	23:DB:1361:G:H5'	1.79	0.65
23:DB:1645:G:H5''	23:DB:1646:C:H5'	1.78	0.65
27:DE:1:MET:HG3	27:DE:18:THR:OG1	1.96	0.65
29:DG:169:ARG:NH1	29:DG:169:ARG:HB2	2.10	0.65
52:DI:108:ILE:HG22	52:DI:128:ILE:HD13	1.79	0.65
36:DO:4:LYS:HD2	36:DO:5:SER:N	2.12	0.65
1:AA:1118:U:H1'	1:AA:1179:A:C4	2.31	0.65
20:AB:112:ARG:O	20:AB:116:LEU:HB2	1.96	0.65
2:AC:19:SER:HB2	2:AC:39:ARG:HH22	1.61	0.65
2:AC:72:PRO:O	2:AC:76:ILE:HG12	1.95	0.65
14:AO:81:ILE:HA	14:AO:86:LEU:HD12	1.78	0.65
23:BB:2133:G:N3	23:BB:2133:G:H2'	2.10	0.65
23:BB:2876:G:H4'	37:BP:2:ASN:HA	1.77	0.65
23:BB:423:A:H5'	23:BB:424:G:C5'	2.26	0.65
23:BB:2204:G:C5'	25:BC:149:LYS:HE3	2.21	0.65
23:BB:2784:U:H5''	26:BD:41:ALA:HB1	1.79	0.65
27:BE:190:ALA:HB1	27:BE:193:VAL:HB	1.79	0.65
36:BO:98:GLN:O	36:BO:103:VAL:HG21	1.97	0.65
41:BT:64:LYS:HB3	41:BT:77:ARG:HH11	1.60	0.65
23:BB:1155:A:H61	45:BY:10:ARG:NH1	1.94	0.65
1:CA:163:C:H2'	1:CA:164:G:O4'	1.97	0.65
5:CF:39:LEU:HD22	5:CF:39:LEU:O	1.97	0.65
47:D0:47:TYR:HB3	47:D0:52:LYS:N	2.11	0.65
23:DB:1856:U:H2'	23:DB:1857:G:O4'	1.96	0.65
23:DB:25:U:H5''	40:DS:80:PRO:HD3	1.78	0.65
52:DI:32:VAL:HG22	52:DI:60:VAL:HG21	1.78	0.65
23:DB:587:C:C3'	33:DL:29:LYS:HD2	2.13	0.65
33:DL:59:ARG:NH1	50:D3:58:ILE:HG21	2.12	0.65
36:DO:25:ARG:HE	36:DO:94:ARG:NH1	1.94	0.65
1:AA:1171:A:H2'	1:AA:1172:C:H6	1.62	0.65
1:AA:312:C:H2'	1:AA:313:A:H8	1.62	0.65
20:AB:185:ILE:HA	20:AB:199:ILE:HB	1.78	0.65
5:AF:17:GLN:HG2	5:AF:21:MET:HG3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1901:A:H1'	25:BC:250:GLN:HE22	1.60	0.65
23:BB:813:U:H2'	23:BB:814:C:C6	2.32	0.65
23:BB:958:U:C4	34:BM:18:ARG:HB2	2.31	0.65
27:BE:128:ALA:HA	27:BE:157:LEU:HD22	1.79	0.65
28:BF:107:VAL:N	28:BF:108:PRO:HD2	2.12	0.65
28:BF:18:GLU:O	28:BF:19:PHE:HB3	1.97	0.65
31:BJ:58:ASN:HA	31:BJ:126:ALA:HA	1.78	0.65
33:BL:63:LYS:HG2	50:B3:11:LYS:HA	1.77	0.65
39:BR:75:VAL:H	39:BR:90:ARG:HE	1.43	0.65
1:CA:1527:U:H2'	1:CA:1528:U:C6	2.31	0.65
1:CA:882:C:O2'	1:CA:883:C:H5'	1.96	0.65
9:CJ:56:HIS:O	9:CJ:57:VAL:HG12	1.96	0.65
10:CK:115:ILE:HD12	10:CK:115:ILE:O	1.96	0.65
12:CM:74:MET:HE2	12:CM:77:LYS:HD2	1.79	0.65
16:CQ:10:ARG:NH2	16:CQ:11:VAL:HB	2.11	0.65
23:DB:179:C:H5''	49:D2:28:ARG:HH12	1.60	0.65
23:DB:222:A:N6	23:DB:232:G:H1'	2.12	0.65
25:DC:171:VAL:HA	25:DC:183:VAL:O	1.97	0.65
28:DF:83:PRO:O	28:DF:84:ILE:HD12	1.96	0.65
29:DG:53:PRO:HG2	29:DG:61:TRP:CZ3	2.32	0.65
32:DK:99:ILE:HD13	32:DK:115:ILE:HG13	1.78	0.65
35:DN:107:ASN:HD22	40:DS:40:ASN:HD22	1.43	0.65
35:DN:42:LYS:HE3	35:DN:42:LYS:HA	1.78	0.65
41:DT:48:GLN:CA	41:DT:53:VAL:HG22	2.25	0.65
1:AA:1471:U:O2'	1:AA:1472:U:H5'	1.97	0.65
1:AA:163:C:H2'	1:AA:164:G:O4'	1.97	0.65
1:AA:882:C:O2'	1:AA:883:C:H5'	1.96	0.65
10:AK:30:ILE:HG22	10:AK:45:THR:HA	1.79	0.65
12:AM:44:ILE:HA	12:AM:47:LEU:HD13	1.79	0.65
14:AO:70:LYS:HZ1	14:AO:74:VAL:HG13	1.62	0.65
19:AT:68:LYS:CG	19:AT:69:ASN:H	2.05	0.65
49:B2:25:LYS:HE3	49:B2:25:LYS:N	2.10	0.65
23:BB:161:A:C3'	23:BB:162:U:H5''	2.19	0.65
23:BB:1979:U:O2'	23:BB:1980:G:H5'	1.96	0.65
27:BE:84:THR:HG23	27:BE:85:PHE:H	1.61	0.65
33:BL:77:ILE:HB	33:BL:110:VAL:CG2	2.26	0.65
1:CA:1069:C:H4'	1:CA:1192:C:O2	1.97	0.65
1:CA:1380:U:O4	6:CG:2:ARG:HB2	1.97	0.65
1:CA:1386:G:H2'	1:CA:1387:G:H8	1.61	0.65
1:CA:372:C:H1'	1:CA:373:A:OP2	1.97	0.65
1:CA:1191:A:OP1	2:CC:3:LYS:HG3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:141:VAL:HG12	3:CD:180:THR:HA	1.77	0.65
5:CF:53:LYS:HD3	5:CF:54:LEU:HB2	1.79	0.65
6:CG:15:PRO:HG2	6:CG:43:TYR:OH	1.96	0.65
8:CI:33:SER:H	8:CI:36:GLN:HB2	1.61	0.65
22:DA:5:U:H2'	22:DA:6:G:C8	2.32	0.65
23:DB:1105:U:H2'	23:DB:1106:G:H8	1.60	0.65
23:DB:1262:A:H2	47:D0:6:LYS:HD2	1.62	0.65
23:DB:139:U:H3'	23:DB:139:U:P	2.37	0.65
23:DB:1515:A:H2'	23:DB:1516:G:O4'	1.97	0.65
23:DB:2537:U:H2'	23:DB:2538:C:H6	1.60	0.65
26:DD:48:ILE:HG22	26:DD:49:GLN:N	2.12	0.65
26:DD:1:MET:HB2	26:DD:81:GLU:OE1	1.97	0.65
30:DH:53:GLU:OE1	30:DH:57:LYS:HD2	1.96	0.65
33:DL:77:ILE:HB	33:DL:110:VAL:HA	1.77	0.65
35:DN:73:ASN:O	35:DN:76:VAL:HG12	1.96	0.65
39:DR:6:GLN:NE2	39:DR:41:ILE:HB	2.12	0.65
23:BB:277:G:H5''	23:BB:278:A:N7	2.12	0.65
23:BB:416:U:H2'	23:BB:417:C:C6	2.31	0.65
25:BC:117:SER:HB3	25:BC:127:ASN:ND2	2.12	0.65
29:BG:157:LYS:HB2	29:BG:159:LYS:HG2	1.78	0.65
31:BJ:17:VAL:HG23	31:BJ:139:VAL:HB	1.79	0.65
38:BQ:16:ILE:HD13	38:BQ:34:ALA:HB2	1.77	0.65
43:BW:57:THR:O	43:BW:58:LEU:HB3	1.96	0.65
46:BZ:32:LEU:HD11	46:BZ:47:LYS:HG3	1.78	0.65
1:CA:1036:A:H2'	1:CA:1037:C:C6	2.31	0.65
1:CA:1270:G:H2'	1:CA:1271:A:C8	2.32	0.65
4:CE:87:VAL:HG13	4:CE:88:HIS:H	1.61	0.65
6:CG:63:VAL:HG12	6:CG:126:ALA:HB1	1.78	0.65
13:CN:25:GLU:O	13:CN:29:ILE:HG13	1.97	0.65
47:D0:29:VAL:HB	47:D0:34:GLY:HA2	1.77	0.65
48:D1:24:LYS:C	48:D1:24:LYS:HZ2	1.99	0.65
23:DB:2285:C:C5	48:D1:7:LYS:HE3	2.32	0.65
23:DB:1368:G:H5''	49:D2:25:LYS:CG	2.23	0.65
25:DC:170:TYR:O	25:DC:171:VAL:HG13	1.96	0.65
26:DD:33:ARG:O	26:DD:34:VAL:HG22	1.96	0.65
27:DE:149:ILE:HD11	27:DE:188:MET:N	2.12	0.65
28:DF:7:TYR:O	28:DF:12:VAL:HG23	1.97	0.65
29:DG:10:VAL:HA	29:DG:14:VAL:HG11	1.77	0.65
34:DM:71:LYS:HZ3	34:DM:91:TYR:HB3	1.61	0.65
35:DN:38:LEU:HB3	35:DN:39:PRO:HD3	1.78	0.65
42:DU:11:ILE:HG22	42:DU:12:VAL:H	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1295:U:H2'	1:AA:1296:C:C6	2.32	0.65
20:AB:141:GLU:O	20:AB:145:ASN:HB2	1.97	0.65
10:AK:22:ILE:HD13	10:AK:95:THR:CG2	2.27	0.65
13:AN:27:LYS:HE2	13:AN:48:GLN:HE22	1.62	0.65
30:BH:3:VAL:HB	30:BH:37:VAL:CB	2.26	0.65
37:BP:38:ARG:HH11	37:BP:38:ARG:HA	1.62	0.65
39:BR:24:LYS:HD3	39:BR:66:HIS:NE2	2.12	0.65
39:BR:65:ALA:HB3	39:BR:99:THR:H	1.61	0.65
4:CE:109:ALA:HB1	4:CE:136:VAL:HG13	1.77	0.65
7:CH:11:THR:HA	7:CH:14:ARG:NH1	2.11	0.65
23:DB:1179:G:H2'	23:DB:1180:U:H6	1.62	0.65
23:DB:144:A:C2	41:DT:3:ARG:NH2	2.65	0.65
23:DB:1979:U:O2'	23:DB:1980:G:H5'	1.96	0.65
23:DB:224:U:O4	23:DB:420:C:H5'	1.97	0.65
23:DB:2286:G:H1	48:D1:23:THR:CG2	2.10	0.65
23:DB:4:U:H2'	23:DB:5:A:H8	1.60	0.65
23:DB:936:A:H2'	23:DB:937:C:C6	2.32	0.65
25:DC:224:MET:CA	25:DC:233:GLY:H	2.09	0.65
31:DJ:41:LYS:HD2	31:DJ:44:TYR:HB3	1.79	0.65
38:DQ:50:ARG:CZ	38:DQ:50:ARG:HA	2.26	0.65
24:DV:25:LYS:HE2	24:DV:41:GLU:CB	2.27	0.65
46:DZ:24:ILE:HD13	46:DZ:24:ILE:N	2.11	0.65
1:AA:190:A:H2'	1:AA:191:G:O4'	1.96	0.64
1:AA:205:A:H2'	1:AA:206:C:H6	1.61	0.64
1:AA:87:C:H2'	1:AA:88:U:H4'	1.76	0.64
1:AA:946:A:H2'	1:AA:947:G:H8	1.60	0.64
1:AA:997:U:O2'	1:AA:998:C:H5'	1.97	0.64
6:AG:44:SER:O	6:AG:48:THR:HG23	1.97	0.64
12:AM:9:PRO:O	12:AM:44:ILE:HG12	1.97	0.64
23:BB:144:A:H2'	23:BB:145:C:C6	2.29	0.64
23:BB:1548:A:H2'	23:BB:1549:A:H8	1.63	0.64
23:BB:2144:G:H22	23:BB:2148:G:H1'	1.62	0.64
23:BB:2365:G:H4'	43:BW:65:LYS:NZ	2.12	0.64
23:BB:2784:U:H2'	23:BB:2785:C:H6	1.61	0.64
23:BB:1113:U:H5''	29:BG:1:SER:CA	2.27	0.64
29:BG:84:LYS:HZ2	29:BG:85:LYS:N	1.95	0.64
31:BJ:39:LYS:NZ	38:BQ:69:ARG:HD2	2.11	0.64
32:BK:17:ARG:HD2	32:BK:18:ARG:H	1.62	0.64
33:BL:111:ILE:H	33:BL:111:ILE:HD13	1.62	0.64
46:BZ:66:ILE:HB	46:BZ:67:PRO:HD3	1.80	0.64
1:CA:1072:G:H2'	1:CA:1073:U:C6	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:312:C:H2'	1:CA:313:A:H8	1.62	0.64
20:CB:112:ARG:O	20:CB:116:LEU:HB2	1.97	0.64
5:CF:1:MET:SD	5:CF:67:PRO:HD3	2.37	0.64
8:CI:98:ARG:HG2	8:CI:103:VAL:HG11	1.78	0.64
23:DB:1820:U:H3	25:DC:197:ALA:CB	2.10	0.64
23:DB:1821:A:H5'	25:DC:155:ARG:HH21	1.61	0.64
23:DB:2051:A:H4'	26:DD:145:SER:HB2	1.80	0.64
23:DB:93:G:H2'	23:DB:94:A:O4'	1.98	0.64
25:DC:179:GLU:CD	25:DC:266:ILE:HA	2.17	0.64
23:DB:2484:G:H1'	34:DM:119:LEU:HD12	1.80	0.64
37:DP:50:ARG:HB2	37:DP:50:ARG:HH11	1.61	0.64
46:DZ:17:SER:H	46:DZ:21:VAL:CG1	2.10	0.64
1:AA:501:C:H2'	1:AA:502:A:C8	2.31	0.64
1:AA:913:A:O2'	1:AA:914:A:O5'	2.13	0.64
20:AB:119:GLN:O	20:AB:124:THR:HG23	1.97	0.64
7:AH:64:TYR:HB3	7:AH:69:ALA:HA	1.79	0.64
8:AI:18:VAL:HG22	8:AI:64:ILE:HG12	1.79	0.64
18:AS:4:LEU:HD13	18:AS:9:PHE:N	2.10	0.64
50:B3:29:ARG:HA	50:B3:33:THR:HG23	1.79	0.64
23:BB:1447:C:H2'	23:BB:1448:G:H8	1.61	0.64
23:BB:1935:G:H1'	23:BB:1964:G:N2	2.12	0.64
23:BB:2814:A:H2'	23:BB:2815:C:C6	2.32	0.64
23:BB:396:G:H4'	46:BZ:28:VAL:CG2	2.25	0.64
26:BD:115:GLY:CA	26:BD:165:MET:HG3	2.27	0.64
26:BD:38:LYS:H	26:BD:38:LYS:HE2	1.62	0.64
27:BE:141:MET:SD	27:BE:143:LEU:HB2	2.37	0.64
30:BH:6:LEU:HD12	30:BH:15:LEU:HA	1.78	0.64
52:BI:109:ALA:HB1	52:BI:124:MET:HG3	1.79	0.64
35:BN:99:LYS:HE2	47:B0:38:LEU:O	1.96	0.64
46:BZ:36:VAL:HA	46:BZ:42:PRO:CB	2.23	0.64
1:CA:190:A:H2'	1:CA:191:G:O4'	1.96	0.64
5:CF:66:ALA:HB1	5:CF:67:PRO:HD2	1.79	0.64
23:DB:1149:G:H2'	23:DB:1150:C:H6	1.61	0.64
23:DB:1179:G:H2'	23:DB:1180:U:C6	2.32	0.64
23:DB:2025:C:H2'	23:DB:2026:U:C6	2.33	0.64
23:DB:2852:G:H2'	23:DB:2853:C:C6	2.32	0.64
23:DB:536:G:C5'	38:DQ:52:ARG:HH22	2.10	0.64
29:DG:10:VAL:N	29:DG:11:PRO:HD3	2.12	0.64
34:DM:41:LEU:O	34:DM:93:VAL:HB	1.97	0.64
1:AA:372:C:H1'	1:AA:373:A:OP2	1.97	0.64
1:AA:376:G:H5''	15:AP:5:ARG:CB	2.26	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:389:A:H3'	1:AA:390:U:H6	1.62	0.64
15:AP:4:ILE:HD12	15:AP:66:THR:O	1.97	0.64
50:B3:26:ALA:HB1	50:B3:29:ARG:CG	2.26	0.64
23:BB:142:A:N3	41:BT:2:ILE:HG23	2.11	0.64
23:BB:1515:A:H2'	23:BB:1516:G:O4'	1.97	0.64
23:BB:2287:A:O2'	23:BB:2288:A:H2'	1.98	0.64
23:BB:6:A:H2'	23:BB:7:G:C8	2.33	0.64
25:BC:65:ASP:OD2	25:BC:101:ARG:HB3	1.98	0.64
26:BD:37:VAL:HG22	26:BD:46:ARG:NE	2.11	0.64
31:BJ:39:LYS:HE2	31:BJ:39:LYS:HA	1.79	0.64
34:BM:96:ILE:HG23	34:BM:97:GLN:N	2.10	0.64
43:BW:48:ALA:HA	43:BW:54:ARG:N	2.10	0.64
46:BZ:32:LEU:HD22	46:BZ:33:ASN:N	2.12	0.64
1:CA:1080:A:H4'	4:CE:20:VAL:HG21	1.78	0.64
1:CA:1244:G:H2'	1:CA:1245:C:C6	2.32	0.64
1:CA:89:U:H2'	1:CA:90:C:C6	2.33	0.64
1:CA:952:U:H2'	1:CA:953:G:H8	1.62	0.64
2:CC:2:GLN:N	2:CC:2:GLN:HE21	1.96	0.64
5:CF:29:ILE:HG22	5:CF:34:GLY:HA3	1.78	0.64
1:CA:1373:G:H5''	6:CG:35:LYS:HB2	1.78	0.64
12:CM:102:LYS:NZ	12:CM:102:LYS:HB2	2.12	0.64
50:D3:4:LYS:HZ2	50:D3:60:CYS:H	1.43	0.64
23:DB:1354:A:H2'	23:DB:1355:G:O4'	1.97	0.64
23:DB:2197:U:O2'	23:DB:2198:A:H2'	1.97	0.64
23:DB:418:C:H2'	23:DB:419:U:C6	2.33	0.64
36:DO:30:ARG:HG2	36:DO:31:THR:H	1.62	0.64
1:AA:1294:G:H2'	1:AA:1295:U:H6	1.60	0.64
1:AA:1513:A:H2'	1:AA:1514:G:C8	2.32	0.64
1:AA:441:A:H61	1:AA:493:A:H62	1.45	0.64
47:B0:52:LYS:NZ	47:B0:52:LYS:HA	2.12	0.64
35:BN:118:ARG:HG2	47:B0:52:LYS:HE2	1.79	0.64
22:BA:20:G:H2'	22:BA:21:G:C8	2.32	0.64
23:BB:1181:U:H2'	23:BB:1182:G:H8	1.62	0.64
23:BB:1387:A:H2'	23:BB:1388:G:H8	1.63	0.64
23:BB:2155:U:H2'	23:BB:2156:G:O4'	1.98	0.64
23:BB:840:C:O2'	23:BB:841:G:H5'	1.97	0.64
25:BC:67:LYS:HG3	25:BC:68:ARG:N	2.12	0.64
30:BH:5:LEU:HD13	30:BH:37:VAL:HG11	1.80	0.64
52:BI:20:SER:O	52:BI:25:PRO:HD2	1.98	0.64
32:BK:5:GLN:O	32:BK:21:CYS:HB3	1.98	0.64
39:BR:47:VAL:HG22	39:BR:48:LYS:HG2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BX:19:LEU:HA	44:BX:22:LEU:HB2	1.78	0.64
1:CA:1001:C:H2'	1:CA:1002:G:C8	2.33	0.64
20:CB:15:PHE:HB3	20:CB:42:LEU:HD11	1.78	0.64
2:CC:2:GLN:H	2:CC:2:GLN:HE21	1.44	0.64
18:CS:41:PRO:HA	18:CS:44:ILE:HG13	1.79	0.64
23:DB:2814:A:H2'	23:DB:2815:C:C6	2.32	0.64
32:DK:61:VAL:HG13	32:DK:87:LEU:HD11	1.80	0.64
1:AA:518:C:H2'	1:AA:530:G:C8	2.31	0.64
7:AH:36:ALA:HA	7:AH:39:LEU:HD12	1.79	0.64
10:AK:95:THR:HG23	10:AK:96:ILE:H	1.62	0.64
48:B1:34:GLU:OE1	48:B1:36:LYS:HE2	1.97	0.64
23:BB:2575:C:H5''	26:BD:149:ASN:H	1.61	0.64
23:BB:28:A:H61	23:BB:512:G:H1'	1.62	0.64
23:BB:639:U:H2'	23:BB:640:C:C6	2.33	0.64
25:BC:227:VAL:HG12	25:BC:228:ASP:OD1	1.98	0.64
26:BD:120:GLY:N	26:BD:123:LYS:HG3	2.13	0.64
28:BF:12:VAL:HG13	28:BF:27:VAL:HG21	1.79	0.64
35:BN:7:GLY:O	35:BN:8:ARG:HB2	1.96	0.64
23:BB:2880:C:H1'	35:BN:92:GLY:O	1.98	0.64
23:BB:855:G:H21	43:BW:23:LYS:CD	2.09	0.64
1:CA:83:C:H4'	1:CA:83:C:OP1	1.98	0.64
1:CA:1103:C:O2	20:CB:105:THR:HG21	1.97	0.64
51:D4:23:ILE:CD1	51:D4:24:ARG:H	2.11	0.64
23:DB:1690:A:H2'	23:DB:1691:C:O4'	1.98	0.64
23:DB:1935:G:H1'	23:DB:1964:G:N2	2.12	0.64
23:DB:86:G:OP1	42:DU:30:SER:HB3	1.97	0.64
25:DC:107:LYS:HE3	25:DC:108:GLY:N	2.05	0.64
25:DC:29:PHE:CD1	25:DC:29:PHE:O	2.50	0.64
25:DC:52:HIS:O	25:DC:53:ILE:HB	1.96	0.64
27:DE:115:GLN:HB3	27:DE:117:ARG:NH1	2.12	0.64
28:DF:56:LEU:HD22	28:DF:88:VAL:HG21	1.80	0.64
30:DH:127:GLU:HA	30:DH:144:VAL:O	1.97	0.64
34:DM:131:VAL:HG22	34:DM:133:LYS:N	2.12	0.64
39:DR:63:VAL:HG22	39:DR:64:VAL:N	2.07	0.64
8:AI:49:GLN:NE2	8:AI:79:ARG:HD2	2.13	0.64
12:AM:13:HIS:HB2	12:AM:16:ILE:HG22	1.79	0.64
12:AM:68:LEU:O	12:AM:72:ILE:HD13	1.97	0.64
18:AS:50:VAL:O	18:AS:56:HIS:HA	1.97	0.64
47:B0:28:SER:HB2	47:B0:34:GLY:N	2.12	0.64
23:BB:13:A:H61	23:BB:525:U:H2'	1.63	0.64
23:BB:196:A:H2'	23:BB:196:A:N3	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:923:G:H1'	43:BW:23:LYS:CD	2.23	0.64
25:BC:171:VAL:HB	25:BC:182:LYS:HB3	1.80	0.64
27:BE:58:LYS:O	27:BE:58:LYS:HD3	1.97	0.64
28:BF:75:GLY:O	28:BF:76:PHE:HB2	1.97	0.64
30:BH:83:LYS:HB3	30:BH:91:PHE:CD1	2.31	0.64
33:BL:109:LYS:HA	33:BL:126:ARG:CA	2.27	0.64
23:BB:1225:G:H5''	39:BR:78:ARG:HH22	1.61	0.64
39:BR:90:ARG:HH11	39:BR:91:GLN:H	1.46	0.64
41:BT:3:ARG:HH22	44:BX:19:LEU:HD22	1.61	0.64
4:CE:137:ARG:HH11	4:CE:137:ARG:HG2	1.63	0.64
11:CL:43:LYS:CB	11:CL:44:PRO:HD2	2.17	0.64
12:CM:14:ALA:HB3	12:CM:40:GLU:HA	1.78	0.64
15:CP:28:ARG:CD	15:CP:29:ASN:H	2.09	0.64
23:DB:2196:C:O2'	23:DB:2197:U:H5'	1.97	0.64
23:DB:2238:G:N3	23:DB:2238:G:H2'	2.13	0.64
23:DB:2365:G:H4'	43:DW:65:LYS:HD2	1.80	0.64
23:DB:321:U:OP2	27:DE:130:LYS:HG3	1.98	0.64
23:DB:64:A:H2'	23:DB:65:U:C6	2.32	0.64
33:DL:39:LYS:NZ	33:DL:39:LYS:HA	2.13	0.64
38:DQ:47:ARG:HH12	38:DQ:50:ARG:HG3	1.62	0.64
44:DX:42:LEU:C	44:DX:43:LEU:HD22	2.17	0.64
2:AC:39:ARG:HG3	2:AC:54:ILE:HG21	1.79	0.64
6:AG:3:ARG:HH11	6:AG:3:ARG:HB3	1.63	0.64
6:AG:4:ARG:HE	6:AG:5:VAL:N	1.94	0.64
10:AK:19:VAL:HG12	10:AK:82:GLU:HB2	1.80	0.64
10:AK:52:ARG:HD2	10:AK:53:GLY:N	2.13	0.64
16:AQ:10:ARG:HA	16:AQ:10:ARG:HE	1.63	0.64
23:BB:1198:U:H2'	23:BB:1199:U:C6	2.32	0.64
23:BB:1459:G:O2'	23:BB:1460:U:H5'	1.98	0.64
23:BB:1813:G:N2	25:BC:51:ARG:HG2	2.12	0.64
23:BB:2104:C:H2'	23:BB:2105:U:C5	2.32	0.64
23:BB:2443:C:H2'	23:BB:2444:G:C8	2.31	0.64
23:BB:796:C:H2'	23:BB:797:G:H8	1.62	0.64
25:BC:250:GLN:HB3	25:BC:252:LYS:HE2	1.78	0.64
26:BD:2:ILE:HG13	26:BD:3:GLY:H	1.62	0.64
26:BD:62:LYS:H	26:BD:62:LYS:HD3	1.62	0.64
29:BG:116:LEU:HD21	29:BG:120:ILE:HB	1.79	0.64
52:BI:85:ILE:HD13	52:BI:137:LEU:HD21	1.78	0.64
46:BZ:36:VAL:HG22	46:BZ:42:PRO:HG3	1.78	0.64
1:CA:1402:C:H2'	1:CA:1403:C:O4'	1.98	0.64
2:CC:149:LYS:HB2	2:CC:168:ARG:HG3	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:28:ASP:CB	3:CD:33:ILE:HD12	2.28	0.64
10:CK:61:ALA:O	10:CK:64:VAL:HG22	1.97	0.64
12:CM:47:LEU:H	12:CM:47:LEU:HD23	1.63	0.64
49:D2:34:ARG:HD2	49:D2:43:THR:OG1	1.97	0.64
22:DA:51:G:H2'	22:DA:52:A:H5''	1.78	0.64
23:DB:639:U:H2'	23:DB:640:C:C6	2.33	0.64
23:DB:796:C:H2'	23:DB:797:G:H8	1.62	0.64
23:DB:813:U:H2'	23:DB:814:C:H6	1.63	0.64
26:DD:128:ARG:HD3	26:DD:128:ARG:N	2.13	0.64
31:DJ:98:GLU:O	31:DJ:102:GLU:HG2	1.98	0.64
31:DJ:64:VAL:HG12	31:DJ:65:THR:N	2.11	0.64
37:DP:59:THR:HG23	37:DP:76:HIS:CD2	2.31	0.64
39:DR:80:ARG:O	39:DR:80:ARG:HG3	1.97	0.64
43:DW:13:ARG:NE	43:DW:13:ARG:H	1.90	0.64
1:AA:1077:G:N2	1:AA:1079:G:H3'	2.13	0.64
1:AA:411:A:O2'	1:AA:412:A:H4'	1.98	0.64
20:AB:198:VAL:HG22	20:AB:200:PRO:HD3	1.79	0.64
8:AI:20:ILE:HG23	8:AI:60:LEU:HD13	1.79	0.64
23:BB:1181:U:H2'	23:BB:1182:G:C8	2.32	0.64
23:BB:1354:A:H2'	23:BB:1355:G:O4'	1.97	0.64
23:BB:1360:G:H2'	23:BB:1361:G:H5'	1.79	0.64
23:BB:2743:U:H3'	23:BB:2744:G:H5''	1.80	0.64
23:BB:674:G:H5''	27:BE:71:GLY:N	2.13	0.64
25:BC:16:VAL:HA	25:BC:17:LYS:HZ3	1.61	0.64
28:BF:151:LEU:HD23	28:BF:153:ILE:HD11	1.80	0.64
31:BJ:24:THR:C	31:BJ:25:LEU:HD22	2.17	0.64
31:BJ:35:ARG:HA	31:BJ:35:ARG:NE	2.13	0.64
41:BT:88:LYS:CG	41:BT:89:GLU:H	2.06	0.64
42:BU:3:LYS:HZ2	42:BU:3:LYS:HB2	1.62	0.64
46:BZ:5:ILE:HB	46:BZ:51:VAL:HG13	1.79	0.64
1:CA:376:G:H5''	15:CP:5:ARG:HD3	1.79	0.64
5:CF:43:GLY:HA2	5:CF:58:HIS:CD2	2.33	0.64
12:CM:89:ARG:NH1	12:CM:101:THR:HG21	2.12	0.64
22:DA:7:G:H5''	36:DO:29:HIS:NE2	2.13	0.64
23:DB:1814:G:C4'	25:DC:51:ARG:HG2	2.28	0.64
23:DB:828:U:H4'	23:DB:831:G:N1	2.13	0.64
25:DC:20:ASN:HB2	25:DC:202:ARG:HH11	1.63	0.64
26:DD:197:THR:HG23	26:DD:198:GLY:N	2.07	0.64
27:DE:108:ILE:HG22	27:DE:180:LEU:HD13	1.79	0.64
29:DG:148:ARG:HA	29:DG:161:VAL:CG1	2.28	0.64
23:DB:1007:C:H5''	31:DJ:37:ARG:NH1	2.13	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DK:35:VAL:HA	32:DK:62:VAL:O	1.98	0.64
23:DB:2334:U:N3	43:DW:74:LYS:HE2	2.10	0.64
46:DZ:48:GLN:HB3	46:DZ:51:VAL:CG2	2.28	0.64
1:AA:1111:A:O2'	1:AA:1112:C:H6	1.81	0.64
1:AA:1278:G:H4'	1:AA:1279:G:O5'	1.98	0.64
1:AA:190:A:O5'	1:AA:190:A:H8	1.80	0.64
1:AA:674:G:H2'	1:AA:675:A:H8	1.63	0.64
6:AG:23:ALA:O	6:AG:26:VAL:HG22	1.98	0.64
9:AJ:7:ARG:HA	9:AJ:75:ASP:HB2	1.78	0.64
9:AJ:65:TYR:HB3	13:AN:95:LEU:HD11	1.80	0.64
16:AQ:59:GLU:H	16:AQ:74:LEU:CD2	2.11	0.64
23:BB:1117:C:H2'	23:BB:1118:C:C6	2.32	0.64
23:BB:1690:A:H2'	23:BB:1691:C:O4'	1.98	0.64
23:BB:1812:U:H1'	25:BC:45:ASN:OD1	1.98	0.64
30:BH:112:LYS:O	30:BH:112:LYS:HD2	1.98	0.64
30:BH:89:LYS:NZ	30:BH:123:ARG:HB3	2.13	0.64
34:BM:109:PRO:HA	34:BM:112:LEU:HD22	1.80	0.64
35:BN:113:ILE:O	35:BN:114:GLU:HB2	1.96	0.64
36:BO:109:ALA:HA	36:BO:112:GLU:OE1	1.98	0.64
42:BU:11:ILE:HB	42:BU:69:VAL:HB	1.78	0.64
42:BU:43:LYS:HD2	42:BU:44:HIS:N	2.13	0.64
45:BY:4:ILE:HG12	45:BY:5:LYS:HG3	1.80	0.64
1:CA:1097:C:H2'	1:CA:1098:C:H6	1.61	0.64
1:CA:1306:A:O2'	12:CM:107:THR:HG21	1.98	0.64
7:CH:102:VAL:HG23	7:CH:125:ILE:HB	1.80	0.64
15:CP:40:ASN:HD21	15:CP:42:ILE:HG12	1.63	0.64
19:CT:52:GLU:O	19:CT:56:ILE:HD13	1.98	0.64
23:DB:1319:C:O2'	23:DB:1320:C:H5'	1.98	0.64
23:DB:1387:A:H2'	23:DB:1388:G:H8	1.63	0.64
23:DB:2820:A:C4'	35:DN:4:ARG:HG3	2.27	0.64
26:DD:140:HIS:O	26:DD:141:ARG:HG2	1.97	0.64
23:DB:2811:G:H5'	26:DD:62:LYS:NZ	2.13	0.64
27:DE:31:VAL:HG21	27:DE:104:ALA:HB2	1.79	0.64
12:CM:78:ARG:HG3	28:DF:111:ARG:HH11	1.62	0.64
33:DL:59:ARG:HA	50:D3:11:LYS:HZ1	1.62	0.64
35:DN:8:ARG:NH1	35:DN:46:ARG:HG3	2.13	0.64
24:DV:9:ARG:HH22	24:DV:16:ALA:HB1	1.61	0.64
1:AA:1488:G:O2'	1:AA:1489:G:H5'	1.98	0.64
20:AB:148:GLY:O	20:AB:151:LYS:HE2	1.98	0.64
20:AB:37:VAL:HG22	20:AB:38:HIS:N	2.13	0.64
3:AD:88:ASN:O	3:AD:92:LEU:HD23	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AH:8:ASP:O	7:AH:12:ARG:HG3	1.98	0.64
9:AJ:25:ILE:HG23	9:AJ:29:ALA:HB3	1.79	0.64
48:B1:13:SER:H	48:B1:50:GLU:HA	1.62	0.64
51:B4:14:CYS:SG	51:B4:25:VAL:HG12	2.38	0.64
23:BB:1188:U:H4'	39:BR:83:TYR:CB	2.27	0.64
23:BB:1973:G:H2'	23:BB:1974:C:C6	2.33	0.64
27:BE:1:MET:O	27:BE:14:VAL:HG22	1.98	0.64
33:BL:132:ARG:O	33:BL:136:GLU:HB3	1.98	0.64
1:CA:1237:C:H3'	1:CA:1238:A:H5'	1.78	0.64
1:CA:451:A:H5'	15:CP:70:ARG:NH2	2.12	0.64
1:CA:677:U:H3	1:CA:713:G:H22	1.47	0.64
1:CA:89:U:H2'	1:CA:90:C:H6	1.62	0.64
1:CA:1526:G:P	21:CU:38:GLU:HB2	2.38	0.64
50:D3:54:LEU:HA	50:D3:57:VAL:HG12	1.80	0.64
23:DB:118:A:OP2	23:DB:119:A:H2'	1.98	0.64
23:DB:1447:C:H2'	23:DB:1448:G:H8	1.61	0.64
23:DB:2249:U:H4'	23:DB:2275:C:C5	2.33	0.64
23:DB:37:C:O2'	23:DB:38:A:H5'	1.98	0.64
23:DB:7:G:H2'	23:DB:8:C:C6	2.33	0.64
23:DB:784:G:H5''	25:DC:225:ASN:HD21	1.63	0.64
25:DC:179:GLU:HG3	25:DC:266:ILE:HG22	1.80	0.64
26:DD:25:THR:HA	26:DD:188:LEU:HD12	1.79	0.64
26:DD:34:VAL:H	26:DD:90:PHE:C	2.01	0.64
30:DH:108:VAL:HG12	30:DH:110:VAL:HB	1.80	0.64
33:DL:78:ARG:NH1	33:DL:78:ARG:HB3	2.13	0.64
38:DQ:91:ARG:HA	38:DQ:94:LEU:HD21	1.79	0.64
43:DW:67:LYS:HD3	43:DW:68:PHE:H	1.62	0.64
1:AA:1412:C:H2'	1:AA:1413:A:H8	1.61	0.63
1:AA:203:G:H1'	1:AA:465:A:N6	2.13	0.63
3:AD:186:GLU:O	3:AD:190:LEU:HD13	1.98	0.63
8:AI:30:ASN:ND2	8:AI:65:THR:HA	2.13	0.63
13:AN:15:LEU:HD12	13:AN:16:ALA:N	2.13	0.63
23:BB:1219:U:H2'	23:BB:1220:G:H8	1.63	0.63
23:BB:704:G:H1'	23:BB:727:A:N6	2.12	0.63
23:BB:1821:A:H5'	25:BC:155:ARG:HH21	1.62	0.63
25:BC:72:GLY:C	25:BC:73:ILE:HD12	2.18	0.63
27:BE:132:LYS:NZ	27:BE:132:LYS:HB2	2.13	0.63
29:BG:23:ILE:O	29:BG:33:THR:HA	1.97	0.63
31:BJ:73:VAL:HG23	31:BJ:74:TYR:H	1.63	0.63
32:BK:108:ARG:HA	32:BK:116:ILE:HG21	1.80	0.63
34:BM:86:LYS:HD2	34:BM:86:LYS:N	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:28:LEU:HD22	35:BN:48:VAL:HG21	1.80	0.63
37:BP:47:ILE:HA	37:BP:63:ILE:HG23	1.79	0.63
40:BS:27:LYS:N	40:BS:27:LYS:HD2	2.09	0.63
1:CA:203:G:H1'	1:CA:465:A:N6	2.13	0.63
1:CA:674:G:H2'	1:CA:675:A:H8	1.63	0.63
10:CK:121:ARG:HE	21:CU:34:ARG:HG2	1.62	0.63
51:D4:10:LEU:HB2	51:D4:25:VAL:HG23	1.80	0.63
23:DB:2080:A:H2'	23:DB:2081:U:C6	2.33	0.63
23:DB:2784:U:H2'	23:DB:2785:C:C6	2.33	0.63
23:DB:878:A:N3	23:DB:878:A:H2'	2.12	0.63
26:DD:146:ILE:H	26:DD:146:ILE:HD12	1.64	0.63
30:DH:1:MET:C	30:DH:21:VAL:HG22	2.19	0.63
34:DM:57:VAL:HG12	34:DM:58:LYS:N	2.09	0.63
43:DW:43:LYS:HD2	43:DW:76:ARG:HA	1.79	0.63
8:AI:122:ARG:HH11	8:AI:122:ARG:HG3	1.64	0.63
48:B1:9:LYS:HD2	48:B1:9:LYS:N	2.14	0.63
50:B3:41:ARG:HD3	50:B3:43:LEU:H	1.63	0.63
22:BA:76:G:H2'	22:BA:77:U:H6	1.63	0.63
23:BB:1192:G:O2'	23:BB:1193:G:H5'	1.98	0.63
23:BB:1205:A:H4'	23:BB:1206:G:OP2	1.96	0.63
23:BB:125:A:H3'	23:BB:126:A:H5'	1.79	0.63
23:BB:298:G:H2'	23:BB:339:U:O4	1.98	0.63
25:BC:141:HIS:CB	25:BC:190:THR:HB	2.29	0.63
26:BD:23:PRO:HB3	26:BD:189:VAL:H	1.63	0.63
43:BW:76:ARG:HG2	43:BW:77:LYS:N	2.13	0.63
1:CA:1084:G:H5'	1:CA:1102:A:OP2	1.98	0.63
1:CA:1404:C:H2'	1:CA:1405:G:C8	2.33	0.63
1:CA:93:U:H2'	1:CA:94:G:H5'	1.79	0.63
23:DB:2108:A:N3	23:DB:2108:A:H2'	2.13	0.63
23:DB:418:C:H2'	23:DB:419:U:H6	1.63	0.63
23:DB:651:G:OP1	50:D3:18:LYS:HG2	1.97	0.63
32:DK:107:LEU:HD21	32:DK:115:ILE:HD13	1.79	0.63
35:DN:30:ARG:NH1	35:DN:74:GLU:HG2	2.14	0.63
36:DO:17:LYS:HD2	36:DO:92:PHE:CD2	2.33	0.63
40:DS:29:VAL:O	40:DS:33:LEU:HD23	1.98	0.63
1:AA:524:G:H2'	1:AA:525:C:C6	2.34	0.63
1:AA:677:U:H3	1:AA:713:G:H22	1.46	0.63
7:AH:11:THR:HA	7:AH:14:ARG:NH1	2.12	0.63
21:AU:7:GLU:OE2	21:AU:15:LEU:HD22	1.98	0.63
23:BB:1060:U:H5	52:BI:131:THR:HG22	1.63	0.63
23:BB:2025:C:H2'	23:BB:2026:U:C6	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2070:A:H2'	23:BB:2071:A:C8	2.33	0.63
25:BC:140:VAL:HG23	25:BC:141:HIS:H	1.64	0.63
30:BH:11:ASN:HD22	30:BH:12:LEU:H	1.46	0.63
30:BH:83:LYS:HB2	30:BH:92:GLY:H	1.63	0.63
35:BN:70:THR:HG21	35:BN:75:ILE:HG22	1.80	0.63
35:BN:75:ILE:HD13	35:BN:75:ILE:O	1.99	0.63
42:BU:35:VAL:HG12	42:BU:36:GLU:N	2.12	0.63
46:BZ:59:ARG:HE	46:BZ:62:LYS:HD2	1.63	0.63
46:BZ:70:LYS:HE2	46:BZ:70:LYS:N	2.12	0.63
1:CA:411:A:O2'	1:CA:412:A:H4'	1.98	0.63
1:CA:441:A:H61	1:CA:493:A:H62	1.45	0.63
1:CA:985:C:H2'	1:CA:986:U:C6	2.33	0.63
8:CI:113:LYS:HA	8:CI:120:ALA:HB2	1.79	0.63
8:CI:86:LEU:HB3	8:CI:93:LEU:HD13	1.80	0.63
11:CL:85:ARG:NH1	11:CL:93:ARG:HB3	2.13	0.63
23:DB:2361:G:H5''	50:D3:27:ASN:HB2	1.79	0.63
50:D3:4:LYS:HE3	50:D3:61:LEU:H	1.63	0.63
23:DB:1442:U:H2'	23:DB:1443:U:C6	2.33	0.63
23:DB:1485:U:H2'	23:DB:1486:U:H6	1.63	0.63
25:DC:75:ALA:HB2	25:DC:95:TYR:CD1	2.33	0.63
27:DE:42:GLY:O	27:DE:43:THR:HG23	1.97	0.63
28:DF:56:LEU:O	28:DF:59:ILE:HG22	1.97	0.63
31:DJ:41:LYS:HG2	38:DQ:63:ARG:NH1	2.12	0.63
23:DB:907:G:H5'	34:DM:24:THR:HG21	1.80	0.63
37:DP:26:GLU:HA	37:DP:47:ILE:N	2.11	0.63
37:DP:27:VAL:HG13	37:DP:47:ILE:CD1	2.26	0.63
37:DP:27:VAL:CG1	37:DP:47:ILE:HD12	2.26	0.63
38:DQ:57:ARG:HH21	38:DQ:92:LYS:HZ3	1.46	0.63
42:DU:28:LEU:HD11	42:DU:31:GLY:C	2.19	0.63
42:DU:13:LEU:HD21	42:DU:69:VAL:HG13	1.79	0.63
6:AG:74:VAL:HG12	6:AG:75:LYS:H	1.62	0.63
11:AL:20:VAL:O	11:AL:23:LEU:HG	1.97	0.63
12:AM:88:LEU:O	12:AM:92:ARG:HG3	1.97	0.63
21:AU:44:ARG:HG3	21:AU:44:ARG:HH11	1.64	0.63
23:BB:1570:A:H2'	23:BB:1571:A:C8	2.34	0.63
23:BB:1641:A:H2'	23:BB:1642:G:O4'	1.99	0.63
23:BB:364:C:H2'	23:BB:365:U:C6	2.33	0.63
23:BB:41:C:H2'	23:BB:42:A:C8	2.33	0.63
23:BB:828:U:H4'	23:BB:831:G:N1	2.13	0.63
26:BD:8:LYS:HB3	37:BP:5:LYS:HZ1	1.61	0.63
27:BE:113:VAL:HG13	27:BE:114:ARG:H	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:BE:132:LYS:HZ2	27:BE:132:LYS:HB2	1.64	0.63
33:BL:90:VAL:CG1	33:BL:121:THR:H	2.10	0.63
34:BM:67:VAL:HB	34:BM:100:LYS:HD3	1.79	0.63
1:CA:390:U:H2'	1:CA:391:G:H8	1.62	0.63
1:CA:947:G:H2'	1:CA:948:C:H6	1.63	0.63
20:CB:163:ILE:HG23	20:CB:164:ASP:N	2.13	0.63
2:CC:53:ARG:HG3	2:CC:113:LYS:HD3	1.80	0.63
8:CI:48:ARG:HA	8:CI:51:LEU:HD12	1.81	0.63
21:CU:34:ARG:HG2	21:CU:35:GLU:N	2.10	0.63
23:DB:1151:A:H2'	23:DB:1152:C:C6	2.33	0.63
23:DB:1278:C:OP1	35:DN:36:THR:HG23	1.98	0.63
23:DB:598:U:H2'	23:DB:599:A:C8	2.34	0.63
23:DB:608:A:H2'	23:DB:609:A:C8	2.34	0.63
26:DD:61:THR:HG23	26:DD:62:LYS:N	2.14	0.63
32:DK:107:LEU:HD21	32:DK:115:ILE:CD1	2.28	0.63
33:DL:78:ARG:HH12	33:DL:80:SER:HB3	1.63	0.63
37:DP:5:LYS:HA	37:DP:5:LYS:HZ3	1.63	0.63
41:DT:21:SER:HA	41:DT:24:MET:SD	2.38	0.63
42:DU:41:VAL:HA	42:DU:57:ILE:HD12	1.80	0.63
1:AA:975:A:H5''	1:AA:976:G:O5'	1.99	0.63
20:AB:71:THR:HG22	20:AB:94:ARG:HH21	1.63	0.63
5:AF:40:GLU:HG3	5:AF:42:TRP:HE1	1.64	0.63
6:AG:65:LEU:O	6:AG:69:ARG:HG3	1.99	0.63
22:BA:28:C:H5'	22:BA:29:A:OP2	1.98	0.63
23:BB:1351:C:H4'	23:BB:1572:A:O4'	1.99	0.63
23:BB:2784:U:H2'	23:BB:2785:C:C6	2.33	0.63
27:BE:3:LEU:O	27:BE:4:VAL:HG12	1.98	0.63
31:BJ:64:VAL:HG12	31:BJ:65:THR:H	1.64	0.63
32:BK:11:ALA:HB3	32:BK:85:VAL:HG22	1.81	0.63
1:CA:1143:G:H2'	1:CA:1144:G:C8	2.34	0.63
1:CA:1323:G:H2'	1:CA:1324:A:C8	2.34	0.63
1:CA:1409:C:H2'	1:CA:1410:A:C8	2.33	0.63
1:CA:720:C:H4'	17:CR:38:ILE:HD11	1.81	0.63
2:CC:149:LYS:HG3	2:CC:166:TRP:HE1	1.63	0.63
7:CH:78:SER:HB2	7:CH:124:ILE:O	1.97	0.63
7:CH:86:LYS:HD2	7:CH:90:GLU:HG2	1.81	0.63
8:CI:54:VAL:HG12	8:CI:93:LEU:HD22	1.80	0.63
23:DB:1205:A:H62	27:DE:164:LEU:HD11	1.63	0.63
23:DB:1854:A:H62	23:DB:1888:G:H8	1.46	0.63
23:DB:85:G:OP1	42:DU:4:ILE:HG23	1.98	0.63
26:DD:42:ASN:O	26:DD:43:ASP:HB2	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:2:GLU:HA	27:DE:16:GLU:HB3	1.81	0.63
34:DM:97:GLN:N	34:DM:98:PRO:CD	2.61	0.63
35:DN:79:LEU:HD22	35:DN:83:LEU:HD23	1.78	0.63
37:DP:23:ASP:O	37:DP:25:VAL:N	2.32	0.63
23:DB:143:C:N3	41:DT:3:ARG:NH1	2.46	0.63
1:AA:105:G:H2'	1:AA:106:C:H6	1.64	0.63
1:AA:1446:A:C2'	1:AA:1447:A:H5''	2.29	0.63
1:AA:74:A:H2'	1:AA:75:G:C5'	2.27	0.63
1:AA:875:U:O2'	7:AH:14:ARG:HD2	1.97	0.63
3:AD:59:LYS:O	3:AD:63:ILE:HG13	1.98	0.63
7:AH:11:THR:HG22	7:AH:15:ASN:ND2	2.13	0.63
16:AQ:10:ARG:NH2	16:AQ:11:VAL:HG23	2.14	0.63
23:BB:1060:U:O4	23:BB:1088:A:N6	2.32	0.63
23:BB:1132:U:C5'	31:BJ:85:LYS:HZ3	2.11	0.63
23:BB:1509:A:H5'	23:BB:1510:G:H5'	1.80	0.63
23:BB:1567:G:H5''	25:BC:84:PRO:HG3	1.81	0.63
23:BB:2498:C:O2'	23:BB:2499:C:H5'	1.98	0.63
23:BB:2852:G:H2'	23:BB:2853:C:C6	2.33	0.63
23:BB:598:U:H2'	23:BB:599:A:C8	2.34	0.63
23:BB:650:C:H5''	50:B3:22:LYS:NZ	2.13	0.63
23:BB:813:U:H2'	23:BB:814:C:H6	1.62	0.63
23:BB:967:U:H2'	23:BB:968:C:C6	2.33	0.63
25:BC:12:ARG:NE	25:BC:18:VAL:HG11	2.14	0.63
29:BG:86:LEU:HD13	29:BG:163:TYR:HA	1.79	0.63
33:BL:38:GLN:O	33:BL:41:ARG:HD3	1.98	0.63
35:BN:37:THR:HA	35:BN:109:PRO:O	1.98	0.63
35:BN:113:ILE:HG22	35:BN:114:GLU:N	2.10	0.63
36:BO:57:ALA:CB	36:BO:62:LEU:HD22	2.18	0.63
37:BP:97:TYR:O	37:BP:98:TYR:HB2	1.97	0.63
24:BV:46:LYS:O	24:BV:50:MET:HG3	1.99	0.63
24:BV:69:GLU:HG2	24:BV:70:ILE:N	2.14	0.63
24:BV:89:ILE:HD12	24:BV:89:ILE:O	1.99	0.63
1:CA:1041:G:H2'	1:CA:1042:A:H8	1.64	0.63
1:CA:1191:A:OP1	2:CC:2:GLN:HB2	1.98	0.63
2:CC:52:SER:O	2:CC:53:ARG:HG3	1.98	0.63
5:CF:79:ARG:NH2	5:CF:87:SER:HB3	2.14	0.63
13:CN:13:VAL:HG22	13:CN:59:GLN:NE2	2.14	0.63
5:CF:86:ARG:NH1	17:CR:63:TYR:HB3	2.13	0.63
50:D3:24:LYS:NZ	50:D3:24:LYS:HB3	2.13	0.63
50:D3:33:THR:C	50:D3:34:LYS:HD2	2.19	0.63
23:DB:1439:A:C6	23:DB:1552:A:N7	2.67	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2498:C:O2'	23:DB:2499:C:H5'	1.99	0.63
23:DB:365:U:H2'	23:DB:366:C:H6	1.61	0.63
23:DB:532:A:H2'	23:DB:532:A:N3	2.13	0.63
23:DB:64:A:H2'	23:DB:65:U:H6	1.64	0.63
26:DD:9:VAL:HG12	26:DD:10:GLY:H	1.62	0.63
31:DJ:105:VAL:HG11	31:DJ:122:LEU:CD1	2.28	0.63
31:DJ:77:HIS:HB2	31:DJ:85:LYS:HE2	1.81	0.63
32:DK:93:GLN:HG2	32:DK:94:PRO:HD2	1.79	0.63
33:DL:126:ARG:HH12	33:DL:128:THR:HG22	1.64	0.63
38:DQ:3:VAL:C	38:DQ:4:LYS:HD3	2.19	0.63
43:DW:67:LYS:HG2	43:DW:71:LYS:CB	2.29	0.63
1:AA:1308:U:H2'	1:AA:1309:G:C8	2.33	0.63
1:AA:10:A:OP2	4:AE:130:THR:HB	1.98	0.63
5:AF:53:LYS:CD	5:AF:54:LEU:H	2.10	0.63
10:AK:58:THR:HG23	10:AK:61:ALA:HB2	1.80	0.63
18:AS:28:LYS:HD3	18:AS:28:LYS:H	1.63	0.63
23:BB:934:U:H2'	23:BB:935:C:H6	1.61	0.63
27:BE:152:GLU:HG2	27:BE:171:ASP:HB3	1.80	0.63
40:BS:47:VAL:HG12	40:BS:103:ILE:HD13	1.80	0.63
43:BW:43:LYS:CG	43:BW:76:ARG:HE	2.12	0.63
1:CA:32:A:H2'	1:CA:33:A:C8	2.34	0.63
20:CB:70:GLY:O	20:CB:92:ASN:HA	1.97	0.63
5:CF:3:HIS:CA	5:CF:65:GLU:HG3	2.28	0.63
23:DB:1351:C:H4'	23:DB:1572:A:O4'	1.99	0.63
23:DB:863:A:H2'	23:DB:864:G:C8	2.34	0.63
25:DC:137:GLY:C	25:DC:139:THR:H	2.02	0.63
28:DF:177:ARG:HE	28:DF:177:ARG:C	2.01	0.63
52:DI:5:GLN:HB2	52:DI:30:GLN:OE1	1.99	0.63
31:DJ:112:GLY:O	31:DJ:116:ARG:HB2	1.99	0.63
33:DL:4:ASN:O	33:DL:5:THR:HG22	1.98	0.63
35:DN:108:ALA:HB1	35:DN:109:PRO:HD2	1.81	0.63
1:AA:1254:A:OP1	9:AJ:47:GLU:HG3	1.99	0.63
1:AA:1524:C:H2'	1:AA:1525:G:H8	1.63	0.63
1:AA:57:G:H2'	1:AA:58:C:H6	1.63	0.63
6:AG:87:PRO:HG2	6:AG:151:ALA:HB2	1.81	0.63
1:AA:982:U:H3'	13:AN:5:MET:HE1	1.81	0.63
13:AN:81:ILE:HD12	13:AN:82:LYS:H	1.64	0.63
14:AO:60:SER:HA	14:AO:63:ARG:HH12	1.64	0.63
47:B0:6:LYS:HA	47:B0:6:LYS:HZ3	1.63	0.63
23:BB:1453:A:H61	35:BN:74:GLU:HG2	1.61	0.63
23:BB:1439:A:C6	23:BB:1552:A:N7	2.67	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:263:G:H2'	23:BB:264:C:O4'	1.98	0.63
23:BB:38:A:O4'	27:BE:46:GLN:CG	2.40	0.63
23:BB:903:C:H2'	23:BB:904:G:H8	1.63	0.63
27:BE:48:THR:HA	27:BE:86:ALA:HB1	1.81	0.63
26:BD:25:THR:HB	37:BP:5:LYS:NZ	2.13	0.63
38:BQ:57:ARG:HA	38:BQ:60:TRP:HE3	1.63	0.63
41:BT:6:ARG:HG3	41:BT:6:ARG:O	1.97	0.63
42:BU:7:ASP:HA	42:BU:24:VAL:HA	1.81	0.63
46:BZ:64:PHE:HA	46:BZ:67:PRO:HD2	1.81	0.63
1:CA:1520:C:H2'	1:CA:1521:C:C6	2.34	0.63
1:CA:976:G:OP1	13:CN:70:HIS:HA	1.98	0.63
2:CC:52:SER:O	2:CC:113:LYS:HD3	1.99	0.63
10:CK:124:LYS:CD	21:CU:34:ARG:HD3	2.28	0.63
49:D2:33:ARG:NE	49:D2:33:ARG:HA	2.13	0.63
23:DB:1570:A:H2'	23:DB:1571:A:C8	2.34	0.63
23:DB:181:A:H2'	23:DB:182:A:C8	2.33	0.63
23:DB:1973:G:H2'	23:DB:1974:C:C6	2.33	0.63
23:DB:2813:A:H2'	23:DB:2814:A:C8	2.34	0.63
23:DB:283:G:H2'	23:DB:284:U:C6	2.34	0.63
25:DC:19:VAL:HG12	25:DC:20:ASN:N	2.13	0.63
27:DE:169:VAL:HG13	27:DE:170:ARG:N	2.08	0.63
1:AA:1102:A:H2'	1:AA:1103:C:C6	2.34	0.63
1:AA:1163:A:H2'	1:AA:1164:G:C8	2.34	0.63
1:AA:1513:A:H2'	1:AA:1514:G:H8	1.63	0.63
1:AA:168:G:O2'	1:AA:169:C:H5'	1.99	0.63
1:AA:810:C:O2'	1:AA:811:C:H5'	1.99	0.63
48:B1:51:ALA:C	48:B1:52:LYS:HD2	2.18	0.63
49:B2:33:ARG:O	49:B2:37:LYS:HB2	1.98	0.63
23:BB:1309:G:H4'	49:B2:7:PRO:HG2	1.80	0.63
23:BB:2212:A:N7	23:BB:2214:C:N4	2.46	0.63
23:BB:2228:G:H2'	23:BB:2229:U:H6	1.63	0.63
23:BB:41:C:H2'	23:BB:42:A:H8	1.63	0.63
23:BB:588:U:OP1	33:BL:28:GLY:N	2.31	0.63
25:BC:131:MET:HE1	25:BC:143:VAL:HG11	1.79	0.63
27:BE:3:LEU:HB2	27:BE:12:LEU:HD22	1.80	0.63
29:BG:17:LYS:HE3	29:BG:24:THR:HG21	1.80	0.63
23:BB:2563:U:H5'	32:BK:27:GLY:HA2	1.79	0.63
32:BK:98:ARG:C	32:BK:99:ILE:HD12	2.20	0.63
34:BM:3:GLN:O	34:BM:6:ARG:N	2.32	0.63
34:BM:96:ILE:HD13	34:BM:97:GLN:HG3	1.80	0.63
37:BP:91:VAL:HB	37:BP:113:LEU:HG	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BS:66:ILE:H	40:BS:66:ILE:HD12	1.64	0.63
1:CA:1234:C:H2'	1:CA:1235:U:C6	2.34	0.63
1:CA:1342:C:H2'	1:CA:1343:G:H8	1.64	0.63
1:CA:57:G:H2'	1:CA:58:C:H6	1.63	0.63
1:CA:763:G:H2'	1:CA:764:C:C6	2.34	0.63
13:CN:19:TYR:HE2	13:CN:50:LEU:HD13	1.64	0.63
50:D3:4:LYS:HG3	50:D3:61:LEU:HB2	1.81	0.63
22:DA:43:C:H4'	28:DF:62:GLN:HE21	1.64	0.63
23:DB:1022:G:C8	31:DJ:68:LYS:HE3	2.34	0.63
23:DB:1174:U:H1'	23:DB:1176:U:C4	2.34	0.63
23:DB:1548:A:H2'	23:DB:1549:A:H8	1.63	0.63
25:DC:16:VAL:HG12	25:DC:16:VAL:O	1.97	0.63
25:DC:19:VAL:HB	25:DC:205:GLY:HA2	1.80	0.63
23:DB:782:A:O2'	25:DC:223:ALA:HB1	1.98	0.63
31:DJ:131:ASN:C	31:DJ:133:ALA:H	2.01	0.63
31:DJ:69:ARG:HH11	31:DJ:69:ARG:HG3	1.62	0.63
23:DB:631:A:O2'	33:DL:66:PHE:HB3	1.98	0.63
35:DN:45:ARG:NH2	35:DN:113:ILE:HD12	2.12	0.63
1:AA:278:G:N2	1:AA:279:A:H62	1.97	0.62
12:AM:22:TYR:HB3	12:AM:69:ARG:HH22	1.63	0.62
23:BB:1158:C:H5''	45:BY:30:ARG:HD3	1.81	0.62
23:BB:1386:C:H2'	23:BB:1387:A:C8	2.34	0.62
23:BB:1442:U:H2'	23:BB:1443:U:C6	2.33	0.62
23:BB:1854:A:H62	23:BB:1888:G:H8	1.46	0.62
23:BB:2415:G:H2'	23:BB:2416:C:H6	1.62	0.62
23:BB:608:A:H2'	23:BB:609:A:C8	2.34	0.62
23:BB:65:U:H2'	23:BB:66:C:H6	1.64	0.62
23:BB:856:G:H2'	23:BB:857:G:C8	2.34	0.62
25:BC:117:SER:HB3	25:BC:127:ASN:CG	2.19	0.62
26:BD:115:GLY:HA3	26:BD:166:GLY:HA2	1.80	0.62
32:BK:99:ILE:HD13	32:BK:118:LEU:HD13	1.80	0.62
34:BM:34:LYS:O	34:BM:124:LEU:HB2	1.99	0.62
37:BP:63:ILE:H	37:BP:63:ILE:HD12	1.64	0.62
42:BU:27:VAL:HA	42:BU:32:LYS:O	1.99	0.62
20:CB:65:LYS:HB2	20:CB:157:PRO:HA	1.80	0.62
2:CC:182:ASP:HB3	2:CC:201:ILE:HB	1.80	0.62
5:CF:38:ARG:HH21	5:CF:63:ASN:HD21	1.46	0.62
16:CQ:11:VAL:HG22	16:CQ:22:VAL:HG22	1.81	0.62
29:DG:168:VAL:HG23	51:D4:31:PRO:HB3	1.81	0.62
23:DB:1548:A:H2'	23:DB:1549:A:C8	2.34	0.62
23:DB:1571:A:H2'	23:DB:1572:A:C8	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1682:G:H2'	23:DB:1683:U:C6	2.34	0.62
23:DB:1722:A:H2'	23:DB:1723:G:C8	2.34	0.62
23:DB:6:A:H2'	23:DB:7:G:C8	2.34	0.62
23:DB:967:U:H2'	23:DB:968:C:C6	2.34	0.62
25:DC:131:MET:HG3	25:DC:187:CYS:SG	2.39	0.62
52:DI:20:SER:O	52:DI:25:PRO:HD2	1.99	0.62
52:DI:45:THR:CA	52:DI:48:ILE:HG22	2.29	0.62
34:DM:18:ARG:CD	34:DM:18:ARG:H	2.09	0.62
37:DP:89:GLY:HA2	37:DP:112:ARG:HH22	1.64	0.62
42:DU:66:VAL:HG22	42:DU:67:SER:N	2.08	0.62
1:AA:1074:G:H2'	1:AA:1075:U:H6	1.64	0.62
10:AK:58:THR:HG23	10:AK:61:ALA:CB	2.30	0.62
16:AQ:10:ARG:HH11	16:AQ:55:GLY:N	1.96	0.62
49:B2:35:ARG:HH12	49:B2:39:ARG:HG2	1.63	0.62
23:BB:1590:A:H2'	23:BB:1591:A:H8	1.64	0.62
23:BB:2574:G:O2'	26:BD:151:THR:HG21	1.99	0.62
23:BB:381:G:O2'	23:BB:382:A:H5'	1.99	0.62
23:BB:444:C:H5''	27:BE:44:ARG:NH1	2.14	0.62
23:BB:547:A:H3'	23:BB:548:G:C8	2.28	0.62
25:BC:245:THR:O	25:BC:247:TRP:N	2.32	0.62
23:BB:2227:A:H4'	25:BC:258:SER:HA	1.80	0.62
28:BF:128:SER:HA	28:BF:154:THR:HG23	1.81	0.62
29:BG:158:GLY:HA3	29:BG:170:THR:HG21	1.80	0.62
39:BR:41:ILE:O	39:BR:43:ASN:N	2.31	0.62
39:BR:78:ARG:HD3	39:BR:88:GLY:N	2.14	0.62
44:BX:43:LEU:HA	44:BX:47:ARG:HB3	1.81	0.62
1:CA:524:G:H2'	1:CA:525:C:C6	2.34	0.62
1:CA:939:G:H5''	6:CG:101:ARG:HH22	1.63	0.62
8:CI:29:ILE:HG23	8:CI:32:ARG:HB2	1.81	0.62
1:CA:1250:A:C4'	8:CI:69:GLY:H	2.10	0.62
23:DB:2742:G:H5''	51:D4:38:GLY:HA2	1.80	0.62
23:DB:1060:U:O4	23:DB:1088:A:N6	2.31	0.62
23:DB:1386:C:H2'	23:DB:1387:A:C8	2.34	0.62
23:DB:2848:G:H1'	23:DB:2868:A:N6	2.14	0.62
25:DC:10:PRO:CD	25:DC:202:ARG:HH22	2.12	0.62
32:DK:43:ILE:HG12	32:DK:52:VAL:HG12	1.81	0.62
32:DK:66:LYS:HA	32:DK:79:PHE:O	1.99	0.62
33:DL:2:ARG:HG3	33:DL:2:ARG:HH11	1.64	0.62
36:DO:91:SER:HB3	36:DO:115:LEU:HD22	1.81	0.62
37:DP:91:VAL:HG23	37:DP:92:ARG:H	1.65	0.62
40:DS:27:LYS:HA	40:DS:70:LYS:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DU:79:ALA:O	42:DU:96:LYS:HB3	1.99	0.62
24:DV:48:MET:HE1	24:DV:85:LYS:HA	1.81	0.62
44:DX:50:VAL:O	44:DX:54:LYS:HB2	1.99	0.62
1:AA:408:A:H3'	1:AA:409:U:H6	1.63	0.62
1:AA:636:U:H2'	1:AA:637:C:C6	2.35	0.62
9:AJ:30:LYS:HB2	9:AJ:36:VAL:HG21	1.81	0.62
12:AM:89:ARG:HB2	12:AM:96:VAL:HG12	1.81	0.62
13:AN:63:CYS:HB3	13:AN:67:GLY:N	2.08	0.62
23:BB:1319:C:O2'	23:BB:1320:C:H5'	1.98	0.62
23:BB:1779:U:H5	23:BB:1784:A:N7	1.97	0.62
23:BB:2274:A:H2'	23:BB:2276:G:OP1	2.00	0.62
23:BB:244:A:H2'	23:BB:245:G:O4'	2.00	0.62
25:BC:224:MET:HA	25:BC:232:GLY:CA	2.28	0.62
25:BC:46:GLY:O	25:BC:47:ARG:HB2	1.98	0.62
23:BB:2575:C:C4'	26:BD:148:GLN:HA	2.28	0.62
26:BD:4:LEU:HD13	26:BD:79:LEU:CD1	2.28	0.62
26:BD:46:ARG:HD2	26:BD:80:TRP:CZ2	2.34	0.62
29:BG:136:ASP:OD1	29:BG:138:GLN:HB3	1.99	0.62
43:BW:42:THR:HA	43:BW:74:LYS:HA	1.80	0.62
44:BX:28:LEU:HB3	44:BX:34:SER:CA	2.24	0.62
3:CD:148:ALA:O	3:CD:151:GLN:HG2	1.99	0.62
7:CH:35:ILE:O	7:CH:38:VAL:HG12	1.99	0.62
23:DB:1024:G:C3'	23:DB:1025:G:H5''	2.28	0.62
23:DB:1100:C:H2'	23:DB:1101:U:H6	1.64	0.62
23:DB:1802:A:H2'	23:DB:1803:A:C8	2.34	0.62
23:DB:182:A:H2'	23:DB:183:C:H6	1.63	0.62
23:DB:2306:C:H3'	23:DB:2307:G:C5'	2.29	0.62
23:DB:2415:G:H2'	23:DB:2416:C:H6	1.62	0.62
25:DC:3:VAL:HG12	25:DC:4:LYS:H	1.64	0.62
26:DD:125:TRP:CD1	26:DD:127:PHE:HB2	2.34	0.62
35:DN:87:PHE:CD1	35:DN:90:ARG:HB2	2.34	0.62
43:DW:35:ILE:CD1	43:DW:35:ILE:H	2.10	0.62
23:DB:1365:A:OP2	46:DZ:9:TYR:HE2	1.81	0.62
1:AA:918:A:H2'	1:AA:919:A:C8	2.34	0.62
8:AI:11:ARG:NH2	8:AI:12:LYS:HB2	2.14	0.62
10:AK:22:ILE:HD13	10:AK:95:THR:HG21	1.81	0.62
10:AK:56:LYS:H	10:AK:56:LYS:HD3	1.64	0.62
15:AP:71:VAL:HA	15:AP:74:LEU:HG	1.79	0.62
23:BB:1097:U:H2'	23:BB:1098:A:H5'	1.81	0.62
23:BB:1131:G:H1'	23:BB:1133:A:N6	2.14	0.62
23:BB:143:C:H2'	23:BB:144:A:N9	2.13	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:172:A:H2'	23:BB:173:A:C8	2.34	0.62
23:BB:932:U:H5'	23:BB:933:A:OP1	1.99	0.62
25:BC:63:ILE:HG22	25:BC:64:VAL:H	1.64	0.62
27:BE:109:LEU:HD23	27:BE:110:SER:N	2.14	0.62
34:BM:45:GLN:NE2	34:BM:119:LEU:HB3	2.14	0.62
41:BT:36:LYS:C	41:BT:36:LYS:HD3	2.20	0.62
44:BX:30:MET:HG3	44:BX:31:GLN:N	2.14	0.62
1:CA:636:U:H2'	1:CA:637:C:C6	2.34	0.62
12:CM:56:ARG:HH11	12:CM:56:ARG:HG3	1.63	0.62
23:DB:1295:C:H2'	23:DB:1296:G:H8	1.64	0.62
23:DB:2088:A:H2'	23:DB:2089:C:H6	1.64	0.62
23:DB:2292:U:H2'	23:DB:2293:G:C8	2.34	0.62
30:DH:65:ALA:HA	30:DH:68:ARG:HB2	1.81	0.62
52:DI:71:LYS:HB3	52:DI:115:ASP:OD2	1.99	0.62
34:DM:16:ARG:HH22	34:DM:72:PRO:HG2	1.65	0.62
36:DO:71:ALA:HB1	36:DO:106:LEU:HB2	1.81	0.62
36:DO:51:ALA:HB3	36:DO:78:VAL:HG22	1.81	0.62
39:DR:39:LEU:H	39:DR:61:ALA:CB	2.11	0.62
43:DW:13:ARG:HE	43:DW:13:ARG:N	1.88	0.62
4:AE:84:VAL:HG22	4:AE:85:LYS:H	1.65	0.62
13:AN:42:ASN:O	13:AN:46:LYS:HG3	2.00	0.62
23:BB:1485:U:H2'	23:BB:1486:U:H6	1.63	0.62
23:BB:1548:A:H2'	23:BB:1549:A:C8	2.34	0.62
23:BB:1818:U:O3'	25:BC:155:ARG:HB2	2.00	0.62
23:BB:19:A:H5''	38:BQ:20:ALA:O	1.99	0.62
23:BB:479:A:O2'	23:BB:480:A:OP2	2.15	0.62
23:BB:969:G:H2'	23:BB:970:U:C6	2.35	0.62
26:BD:62:LYS:HG2	26:BD:63:PRO:CD	2.29	0.62
33:BL:79:LEU:HD12	33:BL:79:LEU:H	1.64	0.62
34:BM:65:ILE:HG22	34:BM:101:VAL:HA	1.82	0.62
1:CA:1071:C:H2'	1:CA:1072:G:C8	2.33	0.62
1:CA:107:G:H2'	1:CA:108:G:H5''	1.81	0.62
1:CA:1329:A:H5''	12:CM:25:GLY:H	1.63	0.62
1:CA:492:C:C2'	1:CA:493:A:H5''	2.28	0.62
1:CA:980:C:H2'	1:CA:981:U:H5'	1.80	0.62
4:CE:95:MET:HA	4:CE:124:ALA:HB2	1.79	0.62
7:CH:19:ALA:HB3	7:CH:21:LYS:HG2	1.80	0.62
16:CQ:24:ILE:HD13	16:CQ:43:LEU:HD13	1.81	0.62
50:D3:4:LYS:HE3	50:D3:61:LEU:HB2	1.81	0.62
23:DB:1098:A:O5'	52:DI:3:LYS:HG2	2.00	0.62
23:DB:1590:A:H2'	23:DB:1591:A:H8	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2233:U:H2'	23:DB:2234:G:C8	2.35	0.62
23:DB:438:G:H2'	23:DB:439:A:C8	2.34	0.62
25:DC:21:PRO:HD2	25:DC:202:ARG:NH1	2.13	0.62
25:DC:235:GLU:HG3	25:DC:236:GLY:H	1.64	0.62
25:DC:28:PRO:HB2	25:DC:79:ARG:HG2	1.80	0.62
28:DF:18:GLU:C	28:DF:20:ASN:H	2.03	0.62
29:DG:55:ASP:CG	29:DG:56:GLY:H	2.02	0.62
29:DG:75:VAL:O	29:DG:79:THR:HG22	1.99	0.62
37:DP:69:VAL:HG13	37:DP:70:GLU:N	2.12	0.62
1:AA:957:U:H3	1:AA:960:U:H5''	1.65	0.62
20:AB:101:THR:HG23	20:AB:102:ASN:H	1.65	0.62
20:AB:75:ALA:O	20:AB:79:VAL:HB	2.00	0.62
3:AD:147:LYS:H	3:AD:147:LYS:HE3	1.65	0.62
23:BB:2066:C:O2'	23:BB:2067:G:H5'	2.00	0.62
23:BB:2678:C:H2'	23:BB:2679:A:C8	2.35	0.62
23:BB:274:C:H2'	23:BB:275:C:O4'	2.00	0.62
23:BB:2848:G:H1'	23:BB:2868:A:N6	2.14	0.62
23:BB:536:G:H4'	38:BQ:56:PHE:CZ	2.35	0.62
25:BC:220:ARG:CZ	25:BC:220:ARG:HA	2.29	0.62
26:BD:107:VAL:HG23	26:BD:204:LYS:O	2.00	0.62
30:BH:82:SER:HB2	30:BH:94:ILE:HD11	1.80	0.62
38:BQ:101:ASP:HB3	38:BQ:104:ALA:HB3	1.81	0.62
40:BS:4:ILE:HG22	40:BS:5:ALA:H	1.64	0.62
1:CA:1143:G:H2'	1:CA:1144:G:H8	1.64	0.62
1:CA:1392:G:O2'	1:CA:1393:U:H5'	2.00	0.62
1:CA:1409:C:H2'	1:CA:1410:A:H8	1.65	0.62
1:CA:384:G:H2'	1:CA:385:C:C6	2.34	0.62
1:CA:810:C:O2'	1:CA:811:C:H5'	1.99	0.62
20:CB:162:VAL:HG11	20:CB:172:ILE:HD11	1.82	0.62
9:CJ:53:ILE:HG22	9:CJ:61:ALA:HB1	1.81	0.62
13:CN:32:ASP:HB2	13:CN:34:ASN:ND2	2.14	0.62
15:CP:20:VAL:HG23	15:CP:34:GLU:O	1.99	0.62
23:DB:1151:A:H2'	23:DB:1152:C:H6	1.63	0.62
23:DB:1459:G:O2'	23:DB:1460:U:H5'	1.98	0.62
23:DB:340:A:H2'	23:DB:341:C:O4'	1.99	0.62
25:DC:68:ARG:NH2	25:DC:190:THR:HG23	2.14	0.62
27:DE:191:ASP:HA	27:DE:194:LYS:HE3	1.81	0.62
30:DH:96:THR:HG23	30:DH:97:ARG:HG2	1.81	0.62
23:DB:1099:G:C8	52:DI:3:LYS:CB	2.83	0.62
23:DB:997:G:H5'	38:DQ:92:LYS:HG3	1.81	0.62
1:AA:1017:U:H2'	1:AA:1018:G:C8	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:973:G:H1'	9:AJ:56:HIS:ND1	2.15	0.62
20:AB:128:LEU:HG	20:AB:132:GLU:HG2	1.82	0.62
3:AD:20:LEU:O	3:AD:21:LYS:HE3	2.00	0.62
9:AJ:24:GLU:OE1	9:AJ:90:LEU:HD22	1.98	0.62
23:BB:1155:A:H61	45:BY:10:ARG:HH12	1.46	0.62
23:BB:1571:A:H2'	23:BB:1572:A:C8	2.34	0.62
23:BB:2813:A:H2'	23:BB:2814:A:C8	2.34	0.62
23:BB:364:C:H2'	23:BB:365:U:H6	1.65	0.62
23:BB:633:A:H2'	23:BB:634:C:H5'	1.81	0.62
25:BC:107:LYS:H	25:BC:194:VAL:HG21	1.63	0.62
23:BB:1843:C:H5''	25:BC:252:LYS:NZ	2.15	0.62
28:BF:33:ILE:HD11	28:BF:90:LEU:HB2	1.81	0.62
36:BO:1:MET:HG2	36:BO:3:LYS:N	2.15	0.62
37:BP:86:LYS:O	37:BP:87:ARG:HG3	1.99	0.62
44:BX:29:ARG:HA	44:BX:34:SER:HB2	1.82	0.62
5:CF:47:LEU:HD21	5:CF:57:ALA:CB	2.30	0.62
47:D0:41:HIS:CD2	47:D0:46:GLY:HA2	2.35	0.62
23:DB:1495:A:H2'	23:DB:1496:A:C8	2.35	0.62
23:DB:2249:U:H4'	23:DB:2275:C:H5	1.65	0.62
23:DB:28:A:N6	23:DB:512:G:H1'	2.14	0.62
23:DB:337:C:H2'	23:DB:338:G:O4'	1.99	0.62
26:DD:17:GLU:HG2	37:DP:80:VAL:HB	1.81	0.62
26:DD:30:GLU:HG2	26:DD:94:GLN:HE22	1.64	0.62
23:DB:1098:A:C3'	52:DI:4:VAL:N	2.62	0.62
32:DK:33:ALA:HB2	32:DK:39:ILE:HD11	1.81	0.62
33:DL:118:THR:HG23	33:DL:119:PRO:CD	2.29	0.62
34:DM:41:LEU:HD23	34:DM:93:VAL:HG11	1.82	0.62
39:DR:6:GLN:NE2	39:DR:6:GLN:N	2.47	0.62
40:DS:7:HIS:CD2	40:DS:10:ALA:HB2	2.35	0.62
41:DT:66:LYS:HZ3	41:DT:66:LYS:C	2.02	0.62
42:DU:46:LYS:HD3	42:DU:53:GLN:HG3	1.82	0.62
42:DU:71:ILE:HG21	42:DU:102:ILE:HD12	1.80	0.62
1:AA:1005:A:H3'	1:AA:1006:G:H8	1.64	0.62
1:AA:1444:U:H2'	1:AA:1445:U:C6	2.35	0.62
1:AA:763:G:H2'	1:AA:764:C:C6	2.33	0.62
1:AA:841:C:H3'	1:AA:843:U:OP2	1.99	0.62
20:AB:147:LEU:O	20:AB:150:ILE:HG22	2.00	0.62
12:AM:48:SER:O	12:AM:52:ILE:HG22	1.98	0.62
1:AA:135:C:O2	15:AP:1:MET:HB2	1.99	0.62
23:BB:1469:A:H2'	23:BB:1470:A:H8	1.62	0.62
23:BB:162:U:O2'	23:BB:163:C:H5'	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:365:U:H2'	23:BB:366:C:C6	2.35	0.62
23:BB:968:C:H2'	23:BB:969:G:H8	1.64	0.62
25:BC:67:LYS:O	25:BC:68:ARG:HB2	1.99	0.62
28:BF:116:LEU:HG	28:BF:176:PHE:HA	1.80	0.62
28:BF:177:ARG:HE	28:BF:178:LYS:N	1.97	0.62
30:BH:89:LYS:C	30:BH:90:LEU:HD12	2.20	0.62
32:BK:35:VAL:HG12	32:BK:69:VAL:HG22	1.82	0.62
33:BL:94:THR:CB	33:BL:103:ILE:HG12	2.29	0.62
37:BP:52:ARG:HH11	37:BP:54:LEU:HB3	1.63	0.62
38:BQ:111:LYS:HB2	39:BR:52:PRO:HB3	1.82	0.62
43:BW:2:HIS:C	43:BW:3:LYS:HD2	2.20	0.62
43:BW:35:ILE:HG23	43:BW:36:ILE:N	2.11	0.62
20:CB:46:VAL:HA	20:CB:49:PHE:CD2	2.35	0.62
2:CC:61:LYS:HB3	2:CC:61:LYS:NZ	2.14	0.62
9:CJ:85:ASP:HB2	9:CJ:89:ARG:NH2	2.14	0.62
23:DB:1178:C:H2'	23:DB:1179:G:H8	1.63	0.62
23:DB:2147:A:H4'	23:DB:2148:G:C8	2.34	0.62
23:DB:2784:U:H2'	23:DB:2785:C:H6	1.62	0.62
26:DD:31:ALA:HB3	26:DD:95:SER:H	1.65	0.62
27:DE:134:LEU:H	27:DE:134:LEU:HD13	1.65	0.62
35:DN:30:ARG:HH12	35:DN:74:GLU:HG2	1.63	0.62
36:DO:82:ALA:C	36:DO:87:ILE:HG21	2.20	0.62
39:DR:69:GLY:HA2	39:DR:96:VAL:HA	1.81	0.62
42:DU:57:ILE:HD13	42:DU:58:VAL:N	2.13	0.62
1:AA:107:G:H2'	1:AA:108:G:H5''	1.81	0.62
20:AB:156:LEU:HD12	20:AB:156:LEU:H	1.63	0.62
3:AD:163:GLN:HB2	3:AD:164:ARG:NH1	2.14	0.62
13:AN:17:ASP:HA	13:AN:21:ALA:CB	2.30	0.62
10:AK:105:ARG:HH21	21:AU:10:PRO:HG3	1.64	0.62
51:B4:2:LYS:NZ	51:B4:2:LYS:HB3	2.15	0.62
23:BB:1159:U:H2'	23:BB:1160:G:H8	1.63	0.62
23:BB:1802:A:H2'	23:BB:1803:A:C8	2.34	0.62
23:BB:2185:U:H2'	23:BB:2186:G:H8	1.64	0.62
23:BB:2678:C:H2'	23:BB:2679:A:H8	1.65	0.62
23:BB:854:C:H2'	23:BB:855:G:H8	1.65	0.62
23:BB:923:G:O2'	23:BB:924:G:H5'	2.00	0.62
26:BD:106:LYS:HZ1	26:BD:208:LYS:HD3	1.65	0.62
26:BD:32:ASN:ND2	26:BD:94:GLN:HG2	2.15	0.62
27:BE:147:LEU:HB3	27:BE:183:PHE:CD2	2.35	0.62
29:BG:71:LEU:HA	29:BG:74:MET:HG3	1.82	0.62
31:BJ:67:ASN:O	31:BJ:70:THR:HG22	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:BL:79:LEU:HD12	33:BL:79:LEU:N	2.15	0.62
34:BM:3:GLN:NE2	34:BM:5:LYS:H	1.98	0.62
38:BQ:60:TRP:O	38:BQ:64:ILE:HG13	2.00	0.62
43:BW:44:PHE:CE2	43:BW:60:ALA:HB3	2.34	0.62
1:CA:1354:U:H2'	1:CA:1355:G:H8	1.63	0.62
1:CA:408:A:H3'	1:CA:409:U:H6	1.63	0.62
6:CG:30:MET:HG3	6:CG:34:LYS:O	2.00	0.62
6:CG:59:GLU:HA	6:CG:62:GLU:OE2	1.99	0.62
8:CI:94:ARG:HB3	8:CI:98:ARG:NE	2.14	0.62
23:DB:1509:A:H5'	23:DB:1510:G:H5'	1.80	0.62
23:DB:1551:A:C3'	23:DB:1552:A:H5''	2.30	0.62
23:DB:1641:A:H2'	23:DB:1642:G:O4'	1.98	0.62
25:DC:139:THR:HA	25:DC:193:GLU:OE1	2.00	0.62
34:DM:26:VAL:HG12	34:DM:27:SER:H	1.64	0.62
35:DN:70:THR:HG22	35:DN:72:ASP:H	1.65	0.62
46:DZ:49:ARG:C	46:DZ:51:VAL:H	2.03	0.62
1:AA:1238:A:C2	1:AA:1241:G:H1'	2.35	0.62
1:AA:313:A:H2'	1:AA:314:C:C6	2.35	0.62
1:AA:384:G:H2'	1:AA:385:C:C6	2.35	0.62
1:AA:492:C:C2'	1:AA:493:A:H5''	2.28	0.62
1:AA:590:U:H2'	1:AA:591:U:C6	2.35	0.62
3:AD:2:ARG:HH12	3:AD:114:ARG:HD2	1.65	0.62
9:AJ:88:MET:HB2	9:AJ:89:ARG:NH1	2.15	0.62
10:AK:56:LYS:O	10:AK:58:THR:HG22	1.99	0.62
14:AO:87:ARG:HH11	14:AO:87:ARG:HA	1.65	0.62
15:AP:46:LYS:CE	15:AP:46:LYS:H	2.12	0.62
15:AP:46:LYS:HE2	15:AP:46:LYS:N	2.15	0.62
16:AQ:26:ARG:HH21	16:AQ:39:ARG:CZ	2.12	0.62
21:AU:36:PHE:HB3	21:AU:40:PRO:CG	2.30	0.62
23:BB:1722:A:N6	23:BB:1738:G:H1'	2.15	0.62
23:BB:2472:G:H2'	23:BB:2475:C:H42	1.64	0.62
23:BB:436:C:O2'	23:BB:437:U:H5'	2.00	0.62
23:BB:438:G:H2'	23:BB:439:A:H8	1.65	0.62
25:BC:179:GLU:CG	25:BC:266:ILE:HA	2.30	0.62
25:BC:179:GLU:HG3	25:BC:266:ILE:HA	1.82	0.62
25:BC:43:ASN:OD1	25:BC:51:ARG:HD3	1.99	0.62
26:BD:48:ILE:HG13	26:BD:80:TRP:NE1	2.15	0.62
29:BG:124:CYS:HB3	29:BG:129:GLU:O	1.99	0.62
34:BM:69:PRO:HB2	34:BM:71:LYS:HZ1	1.65	0.62
40:BS:88:ARG:HD2	40:BS:93:ALA:HB3	1.81	0.62
45:BY:9:THR:CA	45:BY:55:LYS:HE2	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:189:G:OP2	46:BZ:12:ILE:HG12	1.99	0.62
1:CA:1200:C:C4'	1:CA:1201:A:H5'	2.29	0.62
1:CA:168:G:O2'	1:CA:169:C:H5'	1.99	0.62
2:CC:150:VAL:HG22	2:CC:199:VAL:HG12	1.82	0.62
6:CG:11:ILE:HG21	6:CG:27:ASN:HD21	1.65	0.62
6:CG:87:PRO:HG3	6:CG:148:LYS:HA	1.81	0.62
19:CT:3:ILE:O	19:CT:4:LYS:HB2	2.00	0.62
48:D1:15:GLY:HA3	48:D1:47:ILE:HG21	1.82	0.62
49:D2:17:GLY:O	49:D2:19:ARG:N	2.32	0.62
23:DB:1181:U:H2'	23:DB:1182:G:C8	2.35	0.62
23:DB:2425:A:H5'	23:DB:2427:C:O4'	2.00	0.62
23:DB:3:U:H2'	23:DB:4:U:C6	2.34	0.62
23:DB:401:A:H2'	23:DB:402:A:H8	1.65	0.62
23:DB:664:G:H2'	23:DB:665:U:H6	1.65	0.62
23:DB:1819:A:OP1	25:DC:153:LEU:HB2	1.99	0.62
28:DF:65:LEU:HD23	28:DF:87:LYS:HD2	1.80	0.62
29:DG:94:ARG:HB2	29:DG:105:SER:HB2	1.82	0.62
30:DH:4:ILE:HD11	30:DH:37:VAL:HA	1.81	0.62
34:DM:133:LYS:HD2	34:DM:134:THR:O	2.00	0.62
42:DU:9:GLU:HB2	42:DU:71:ILE:HG13	1.82	0.62
1:AA:1182:G:H4'	1:AA:1183:U:H5'	1.81	0.61
11:AL:43:LYS:N	11:AL:44:PRO:HD2	2.15	0.61
1:AA:275:G:O5'	16:AQ:15:LYS:HG2	2.00	0.61
19:AT:70:LYS:NZ	19:AT:70:LYS:HB2	2.13	0.61
47:B0:41:HIS:HD2	47:B0:42:ILE:HG22	1.65	0.61
23:BB:1117:C:H2'	23:BB:1118:C:H6	1.65	0.61
23:BB:1295:C:H2'	23:BB:1296:G:H8	1.65	0.61
23:BB:1551:A:C3'	23:BB:1552:A:H5''	2.30	0.61
23:BB:2743:U:H2'	23:BB:2744:G:O4'	2.00	0.61
25:BC:103:ILE:HG22	25:BC:104:LEU:N	2.15	0.61
26:BD:23:PRO:HB3	26:BD:189:VAL:N	2.15	0.61
33:BL:61:LEU:HD22	33:BL:61:LEU:H	1.64	0.61
42:BU:5:ARG:CG	42:BU:6:ARG:H	2.13	0.61
43:BW:16:GLU:CG	43:BW:37:VAL:HG22	2.30	0.61
1:CA:590:U:H2'	1:CA:591:U:C6	2.35	0.61
2:CC:18:ASN:O	2:CC:55:VAL:HA	2.00	0.61
6:CG:125:ASP:HB3	6:CG:129:ASN:HA	1.83	0.61
7:CH:6:ILE:HD11	7:CH:31:LEU:HD23	1.82	0.61
1:CA:33:A:H1'	11:CL:27:PRO:HG3	1.81	0.61
48:D1:19:PHE:HD1	48:D1:20:TYR:H	1.48	0.61
23:DB:1164:C:H2'	23:DB:1165:A:C8	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1254:A:N1	27:DE:77:ILE:HD12	2.14	0.61
23:DB:1439:A:N7	23:DB:1440:U:C6	2.68	0.61
23:DB:1779:U:H5	23:DB:1784:A:N7	1.97	0.61
23:DB:2008:C:H2'	23:DB:2009:A:C8	2.31	0.61
23:DB:2405:G:H5'	33:DL:70:LYS:HG3	1.81	0.61
23:DB:2678:C:H2'	23:DB:2679:A:C8	2.35	0.61
23:DB:30:G:H2'	23:DB:31:C:C6	2.34	0.61
23:DB:925:A:O2'	23:DB:926:G:H5'	2.00	0.61
25:DC:210:ALA:HA	25:DC:213:ARG:HD2	1.82	0.61
28:DF:177:ARG:NH2	28:DF:178:LYS:HA	2.14	0.61
34:DM:96:ILE:HD12	34:DM:96:ILE:N	2.15	0.61
36:DO:15:ARG:HH12	43:DW:76:ARG:HE	1.45	0.61
38:DQ:36:GLN:O	38:DQ:39:ILE:HG23	1.99	0.61
43:DW:36:ILE:HG13	43:DW:38:ARG:H	1.65	0.61
1:AA:32:A:H2'	1:AA:33:A:C8	2.34	0.61
20:AB:87:ASP:O	20:AB:88:GLN:HG3	2.01	0.61
1:AA:16:A:O2'	4:AE:20:VAL:HG13	2.00	0.61
7:AH:111:THR:H	7:AH:114:ALA:HB3	1.64	0.61
7:AH:8:ASP:OD2	7:AH:12:ARG:HD2	1.99	0.61
8:AI:80:HIS:O	8:AI:84:ARG:HB2	1.99	0.61
16:AQ:10:ARG:HA	16:AQ:10:ARG:NE	2.15	0.61
23:BB:1439:A:N7	23:BB:1440:U:C6	2.68	0.61
23:BB:1682:G:H2'	23:BB:1683:U:C6	2.35	0.61
23:BB:1722:A:H2'	23:BB:1723:G:C8	2.34	0.61
25:BC:70:LYS:HE3	25:BC:95:TYR:CE2	2.35	0.61
26:BD:22:ILE:N	26:BD:23:PRO:CD	2.63	0.61
26:BD:33:ARG:HG3	26:BD:34:VAL:H	1.64	0.61
26:BD:60:VAL:HB	26:BD:63:PRO:HG2	1.80	0.61
27:BE:105:LEU:HD22	27:BE:177:PRO:HG2	1.81	0.61
27:BE:57:LYS:HA	27:BE:57:LYS:HE3	1.81	0.61
29:BG:86:LEU:HD12	29:BG:161:VAL:HG12	1.82	0.61
29:BG:18:ILE:HA	29:BG:23:ILE:HD13	1.82	0.61
52:BI:135:MET:HG3	52:BI:137:LEU:HG	1.82	0.61
41:BT:13:ALA:O	41:BT:32:LEU:HA	2.00	0.61
24:BV:17:SER:HB3	24:BV:21:ARG:HH12	1.65	0.61
1:CA:1437:A:H2'	1:CA:1438:G:H8	1.65	0.61
1:CA:60:A:H4'	1:CA:61:G:O5'	2.00	0.61
1:CA:859:G:H2'	1:CA:860:A:H8	1.65	0.61
7:CH:124:ILE:HD12	7:CH:125:ILE:N	2.15	0.61
12:CM:6:ILE:O	12:CM:8:ILE:HG23	2.00	0.61
15:CP:68:SER:OG	15:CP:71:VAL:HG12	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:CU:17:ARG:H	21:CU:17:ARG:CD	2.09	0.61
23:DB:1797:G:O3'	25:DC:253:GLY:HA2	1.99	0.61
23:DB:2324:U:H3'	23:DB:2325:G:H5''	1.82	0.61
23:DB:222:A:H61	23:DB:232:G:H1'	1.62	0.61
23:DB:2755:C:H2'	51:D4:19:ARG:HH21	1.64	0.61
23:DB:448:U:H3'	27:DE:79:ARG:HE	1.63	0.61
23:DB:620:G:H5'	23:DB:620:G:N3	2.15	0.61
26:DD:27:ILE:HG12	26:DD:185:ASN:O	2.00	0.61
27:DE:77:ILE:O	27:DE:77:ILE:HG12	1.99	0.61
28:DF:67:THR:OG1	28:DF:85:GLY:HA3	1.99	0.61
29:DG:175:LYS:HG2	29:DG:176:LYS:N	2.14	0.61
52:DI:121:ILE:N	52:DI:121:ILE:HD13	2.15	0.61
23:DB:1064:C:H4'	52:DI:90:GLY:CA	2.28	0.61
31:DJ:45:THR:HG22	31:DJ:47:HIS:H	1.65	0.61
31:DJ:96:ARG:HG3	31:DJ:98:GLU:OE1	1.99	0.61
33:DL:78:ARG:NE	33:DL:113:ALA:HB1	2.14	0.61
23:DB:1225:G:OP1	39:DR:90:ARG:HD2	2.00	0.61
42:DU:9:GLU:HB2	42:DU:71:ILE:CB	2.30	0.61
44:DX:44:LYS:HG2	44:DX:47:ARG:HB2	1.82	0.61
1:AA:1506:U:H4'	10:AK:128:VAL:OXT	2.00	0.61
5:AF:4:TYR:O	5:AF:63:ASN:HA	2.00	0.61
9:AJ:31:ARG:HH11	9:AJ:31:ARG:HG2	1.65	0.61
9:AJ:53:ILE:HG12	9:AJ:63:ASP:HB2	1.82	0.61
10:AK:36:ARG:HG2	10:AK:37:GLN:H	1.65	0.61
14:AO:55:LEU:O	14:AO:59:VAL:HG23	2.00	0.61
48:B1:13:SER:N	48:B1:50:GLU:HA	2.15	0.61
23:BB:1405:U:H2'	23:BB:1406:U:C6	2.36	0.61
23:BB:1791:A:C2	23:BB:1829:A:H4'	2.35	0.61
23:BB:1915:U:H3'	23:BB:1916:A:H8	1.66	0.61
23:BB:198:C:O2'	23:BB:199:A:H5''	2.00	0.61
23:BB:2455:G:H2'	23:BB:2456:C:C6	2.36	0.61
23:BB:526:A:N6	23:BB:2626:C:H4'	2.16	0.61
23:BB:972:A:C3'	23:BB:973:A:H5''	2.30	0.61
25:BC:37:SER:OG	25:BC:62:ARG:HG3	2.00	0.61
26:BD:128:ARG:NH2	26:BD:130:GLN:HG3	2.14	0.61
26:BD:16:THR:C	37:BP:80:VAL:HB	2.20	0.61
35:BN:1:MET:HE3	35:BN:3:HIS:H	1.64	0.61
36:BO:49:VAL:HG22	36:BO:50:ALA:N	2.13	0.61
42:BU:40:LEU:O	42:BU:42:LYS:HD2	2.00	0.61
1:CA:1030:U:H5''	1:CA:1031:C:C5	2.34	0.61
1:CA:1078:U:H4'	4:CE:137:ARG:HH12	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1151:A:O2'	1:CA:1152:A:H5''	2.00	0.61
1:CA:1287:A:H2'	1:CA:1288:A:C8	2.35	0.61
1:CA:841:C:H3'	1:CA:843:U:OP2	1.99	0.61
1:CA:673:A:H4'	5:CF:86:ARG:HD2	1.81	0.61
6:CG:144:ALA:C	6:CG:146:ALA:H	2.04	0.61
21:CU:24:LYS:CD	21:CU:25:ALA:H	2.11	0.61
23:DB:1099:G:H8	52:DI:3:LYS:CB	2.13	0.61
23:DB:1722:A:N6	23:DB:1738:G:H1'	2.15	0.61
23:DB:453:A:H4'	23:DB:472:A:N6	2.15	0.61
23:DB:784:G:O2'	23:DB:785:G:H5''	2.01	0.61
23:DB:912:C:O2'	23:DB:913:U:H5'	1.99	0.61
25:DC:20:ASN:CG	25:DC:202:ARG:HD3	2.20	0.61
33:DL:18:ARG:HH22	33:DL:21:ARG:HD3	1.65	0.61
33:DL:62:PRO:HB3	50:D3:12:ARG:NE	2.16	0.61
23:DB:2293:G:H5''	36:DO:9:ARG:HH21	1.64	0.61
23:DB:144:A:C6	41:DT:3:ARG:NH1	2.68	0.61
43:DW:67:LYS:HD2	43:DW:70:VAL:N	2.15	0.61
1:AA:1074:G:H2'	1:AA:1075:U:C6	2.35	0.61
1:AA:642:A:H2'	1:AA:643:C:C6	2.35	0.61
50:B3:2:LYS:O	50:B3:2:LYS:HD3	2.00	0.61
23:BB:1105:U:H2'	23:BB:1106:G:C8	2.34	0.61
23:BB:1791:A:OP1	25:BC:211:ARG:HG2	2.00	0.61
23:BB:18:U:H2'	23:BB:19:A:H8	1.64	0.61
23:BB:2809:A:H2'	23:BB:2810:A:C8	2.36	0.61
26:BD:128:ARG:CA	26:BD:128:ARG:HE	2.13	0.61
23:BB:2575:C:H4'	26:BD:148:GLN:C	2.20	0.61
26:BD:170:VAL:HG11	26:BD:194:PRO:CB	2.29	0.61
34:BM:42:THR:CG2	34:BM:45:GLN:HB2	2.30	0.61
39:BR:53:PHE:CE2	39:BR:54:VAL:HB	2.36	0.61
43:BW:54:ARG:HG2	43:BW:54:ARG:HH11	1.65	0.61
43:BW:66:VAL:HG22	43:BW:67:LYS:H	1.64	0.61
1:CA:105:G:H2'	1:CA:106:C:H6	1.64	0.61
20:CB:10:LYS:HB2	20:CB:211:LEU:CD2	2.30	0.61
20:CB:23:ASN:HD22	20:CB:24:PRO:HD2	1.63	0.61
22:DA:65:U:O2'	22:DA:66:A:H5'	1.99	0.61
23:DB:1201:U:H2'	23:DB:1202:G:C8	2.35	0.61
23:DB:358:U:H2'	23:DB:359:G:C8	2.35	0.61
23:DB:876:C:C2	23:DB:877:A:H1'	2.35	0.61
23:DB:937:C:H2'	23:DB:938:G:H8	1.65	0.61
25:DC:136:VAL:HA	25:DC:165:ALA:HA	1.82	0.61
30:DH:131:SER:HB2	30:DH:141:LYS:HA	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:DH:3:VAL:HB	30:DH:37:VAL:CG1	2.31	0.61
30:DH:51:ARG:HG2	30:DH:51:ARG:HH11	1.65	0.61
34:DM:28:PHE:HB2	34:DM:102:LEU:HG	1.82	0.61
38:DQ:25:GLY:HA2	38:DQ:29:ARG:HD2	1.82	0.61
39:DR:65:ALA:CB	39:DR:99:THR:HG23	2.25	0.61
39:DR:71:LYS:HB2	39:DR:73:LYS:HZ1	1.64	0.61
40:DS:21:ALA:HB1	40:DS:74:ILE:HD12	1.82	0.61
24:DV:53:LYS:HB3	24:DV:53:LYS:HZ2	1.66	0.61
43:DW:42:THR:HB	43:DW:75:ASN:CG	2.21	0.61
1:AA:1040:U:H2'	1:AA:1041:G:H8	1.66	0.61
8:AI:4:GLN:HG2	8:AI:21:LYS:HG3	1.81	0.61
10:AK:18:GLY:O	10:AK:81:LEU:HB2	2.00	0.61
33:BL:62:PRO:C	50:B3:24:LYS:HD3	2.20	0.61
22:BA:104:A:H2'	22:BA:105:G:O4'	2.00	0.61
23:BB:1439:A:C4	23:BB:1552:A:N6	2.69	0.61
23:BB:1712:U:H3'	23:BB:1713:A:H5''	1.83	0.61
23:BB:1838:C:N4	23:BB:1898:U:H2'	2.15	0.61
23:BB:37:C:O2'	27:BE:46:GLN:HB3	2.00	0.61
25:BC:15:VAL:HG12	25:BC:16:VAL:HG23	1.80	0.61
25:BC:220:ARG:NE	25:BC:220:ARG:HA	2.15	0.61
26:BD:128:ARG:HH22	26:BD:130:GLN:CG	2.13	0.61
23:BB:1656:C:H5''	26:BD:141:ARG:HB3	1.83	0.61
27:BE:130:LYS:HD3	27:BE:132:LYS:HD2	1.82	0.61
28:BF:108:PRO:HB3	28:BF:113:PHE:CE2	2.36	0.61
30:BH:132:PHE:HB3	30:BH:140:ALA:HB3	1.82	0.61
23:BB:831:G:O3'	33:BL:44:GLY:HA3	2.00	0.61
22:BA:27:C:H5'	36:BO:34:HIS:CE1	2.35	0.61
45:BY:16:LEU:N	45:BY:16:LEU:HD12	2.16	0.61
1:CA:1154:G:O2'	1:CA:1155:A:H5'	2.01	0.61
1:CA:974:A:C4'	1:CA:975:A:H5'	2.18	0.61
3:CD:130:ASN:H	3:CD:130:ASN:HD22	1.49	0.61
11:CL:64:SER:OG	11:CL:96:THR:HG23	2.01	0.61
23:DB:1405:U:H2'	23:DB:1406:U:C6	2.36	0.61
23:DB:2678:C:H2'	23:DB:2679:A:H8	1.64	0.61
23:DB:361:G:H2'	23:DB:362:A:H8	1.64	0.61
23:DB:1656:C:OP1	26:DD:141:ARG:HD2	2.00	0.61
23:DB:2574:G:N2	26:DD:147:GLY:HA3	2.11	0.61
29:DG:30:GLY:H	29:DG:78:VAL:HG23	1.65	0.61
30:DH:11:ASN:HD22	30:DH:20:ASN:ND2	1.97	0.61
23:DB:136:G:H21	41:DT:3:ARG:HG2	1.66	0.61
24:DV:48:MET:SD	24:DV:86:LEU:HD12	2.40	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DY:15:ARG:NE	45:DY:15:ARG:HA	2.16	0.61
1:AA:1011:C:H2'	1:AA:1012:A:H8	1.65	0.61
1:AA:1132:C:H2'	1:AA:1133:G:O4'	2.00	0.61
1:AA:1182:G:H4'	1:AA:1183:U:C5'	2.30	0.61
1:AA:1217:C:OP1	13:AN:8:ARG:HD3	2.01	0.61
1:AA:1320:C:C5	18:AS:36:ARG:HG2	2.36	0.61
1:AA:22:G:H4'	1:AA:885:G:C8	2.36	0.61
4:AE:33:THR:O	4:AE:34:ALA:HB3	2.00	0.61
5:AF:6:ILE:HD13	5:AF:89:VAL:HB	1.83	0.61
23:BB:1105:U:H2'	23:BB:1106:G:H8	1.65	0.61
23:BB:2425:A:H5'	23:BB:2427:C:O4'	2.00	0.61
26:BD:129:THR:O	26:BD:140:HIS:HA	2.01	0.61
26:BD:23:PRO:HG2	26:BD:178:VAL:HG21	1.81	0.61
26:BD:48:ILE:HG13	26:BD:80:TRP:HE1	1.64	0.61
27:BE:147:LEU:HD13	27:BE:183:PHE:HA	1.83	0.61
28:BF:126:ASN:H	28:BF:126:ASN:ND2	1.98	0.61
33:BL:126:ARG:NH2	33:BL:131:ALA:HB2	2.16	0.61
37:BP:38:ARG:HG3	37:BP:39:LEU:N	2.15	0.61
38:BQ:39:ILE:HG23	39:BR:80:ARG:HD3	1.82	0.61
39:BR:84:ARG:NH1	39:BR:86:GLN:HB3	2.15	0.61
1:CA:1023:U:H2'	1:CA:1024:G:C8	2.35	0.61
12:CM:9:PRO:HB2	12:CM:17:ALA:HB1	1.83	0.61
23:DB:104:A:H2'	23:DB:105:C:H6	1.66	0.61
23:DB:544:C:H4'	23:DB:545:U:OP1	2.00	0.61
33:DL:110:VAL:HG22	33:DL:127:VAL:HA	1.82	0.61
33:DL:34:GLY:HA3	39:DR:85:LYS:HE2	1.82	0.61
23:DB:2393:U:H4'	33:DL:62:PRO:O	2.01	0.61
23:DB:64:A:C5'	41:DT:76:ARG:HD2	2.30	0.61
42:DU:4:ILE:HG22	42:DU:27:VAL:HG22	1.81	0.61
7:AH:82:LEU:O	7:AH:82:LEU:HD13	1.99	0.61
19:AT:81:GLN:O	19:AT:85:LEU:HB3	2.01	0.61
10:AK:111:ASP:HB2	21:AU:16:ARG:NH2	2.15	0.61
23:BB:1015:U:H2'	23:BB:1016:G:C8	2.36	0.61
23:BB:527:C:H5''	23:BB:2779:U:O2	2.00	0.61
23:BB:322:A:H5'	23:BB:340:A:H1'	1.83	0.61
23:BB:664:G:H2'	23:BB:665:U:H6	1.65	0.61
26:BD:187:LEU:HD23	26:BD:188:LEU:N	2.16	0.61
26:BD:175:LEU:HD22	26:BD:189:VAL:HG12	1.81	0.61
31:BJ:73:VAL:HB	31:BJ:75:TYR:CE1	2.35	0.61
38:BQ:95:ALA:C	38:BQ:97:ILE:H	2.03	0.61
41:BT:65:GLY:O	41:BT:66:LYS:HB3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BV:21:ARG:HA	24:BV:25:LYS:O	2.00	0.61
43:BW:49:ASN:HB2	43:BW:53:GLY:HA3	1.82	0.61
1:CA:1014:A:H4'	18:CS:13:HIS:CD2	2.36	0.61
1:CA:1292:G:H2'	1:CA:1293:C:C6	2.35	0.61
22:DA:25:U:H4'	22:DA:27:C:OP1	2.00	0.61
23:DB:2471:A:O2'	23:DB:2472:G:O5'	2.19	0.61
23:DB:381:G:O2'	23:DB:382:A:H5'	2.00	0.61
23:DB:918:A:H2'	23:DB:919:U:C5'	2.27	0.61
25:DC:20:ASN:CB	25:DC:202:ARG:HD3	2.31	0.61
25:DC:3:VAL:HG21	25:DC:9:SER:HB3	1.82	0.61
28:DF:140:ILE:H	28:DF:140:ILE:CD1	2.13	0.61
52:DI:27:LEU:HB2	52:DI:32:VAL:HG21	1.82	0.61
31:DJ:73:VAL:HG22	31:DJ:74:TYR:N	2.16	0.61
33:DL:77:ILE:HD11	33:DL:109:LYS:HB2	1.83	0.61
26:DD:14:ILE:HD12	37:DP:78:PRO:CG	2.30	0.61
38:DQ:10:ARG:O	38:DQ:14:LYS:HB2	2.01	0.61
39:DR:69:GLY:HA2	39:DR:97:LYS:H	1.66	0.61
40:DS:36:LEU:HA	40:DS:39:THR:HG22	1.81	0.61
45:DY:29:ARG:O	45:DY:33:HIS:HB3	2.00	0.61
1:AA:132:C:H5''	19:AT:68:LYS:HE3	1.83	0.61
11:AL:3:VAL:O	11:AL:7:VAL:HG23	1.99	0.61
12:AM:22:TYR:HB3	12:AM:69:ARG:NH2	2.15	0.61
23:BB:1874:C:H2'	23:BB:1875:G:O4'	2.01	0.61
23:BB:2291:U:H2'	23:BB:2292:U:C6	2.35	0.61
23:BB:2888:C:H2'	23:BB:2889:C:C6	2.36	0.61
23:BB:355:U:H2'	23:BB:356:G:C8	2.35	0.61
23:BB:990:A:H1'	23:BB:1156:A:H2	1.65	0.61
25:BC:23:LEU:O	25:BC:23:LEU:HD12	2.01	0.61
29:BG:5:LYS:HG2	29:BG:65:GLY:HA2	1.83	0.61
52:BI:27:LEU:CD2	52:BI:27:LEU:H	2.11	0.61
33:BL:38:GLN:NE2	33:BL:39:LYS:H	1.98	0.61
23:BB:2423:U:H3	43:BW:1:ALA:HB2	1.66	0.61
43:BW:23:LYS:NZ	43:BW:24:ARG:HG2	2.16	0.61
1:CA:1031:C:H4'	1:CA:1032:G:C5'	2.30	0.61
1:CA:1342:C:H2'	1:CA:1343:G:C8	2.35	0.61
1:CA:313:A:H2'	1:CA:314:C:C6	2.35	0.61
17:CR:20:ILE:O	17:CR:21:ASP:HB2	2.00	0.61
23:DB:1229:C:H2'	23:DB:1230:A:C8	2.36	0.61
23:DB:1821:A:H2'	23:DB:1822:C:C6	2.36	0.61
23:DB:1838:C:N4	23:DB:1898:U:H2'	2.15	0.61
23:DB:2472:G:H2'	23:DB:2475:C:H42	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:285:G:H2'	23:DB:286:U:O4'	2.01	0.61
23:DB:765:C:H2'	23:DB:766:U:H6	1.66	0.61
25:DC:74:PRO:HG3	25:DC:116:GLN:HE21	1.64	0.61
25:DC:257:ARG:HA	25:DC:261:ARG:CD	2.30	0.61
26:DD:48:ILE:HA	26:DD:79:LEU:O	2.01	0.61
29:DG:51:PHE:HE1	29:DG:53:PRO:HG3	1.66	0.61
30:DH:122:LEU:HA	30:DH:146:VAL:HG21	1.83	0.61
30:DH:66:ASN:HA	30:DH:138:VAL:HG22	1.83	0.61
31:DJ:99:ARG:O	31:DJ:103:ILE:HG12	2.00	0.61
33:DL:110:VAL:O	33:DL:111:ILE:HG12	2.00	0.61
45:DY:2:LYS:HD2	45:DY:35:VAL:HB	1.82	0.61
46:DZ:62:LYS:HA	46:DZ:65:ASN:HD21	1.66	0.61
1:AA:72:A:H2'	1:AA:73:C:C6	2.36	0.61
10:AK:97:ARG:HH11	10:AK:97:ARG:HB3	1.66	0.61
14:AO:2:LEU:HD23	14:AO:3:SER:N	2.16	0.61
23:BB:1019:U:H2'	23:BB:1020:A:C8	2.36	0.61
23:BB:2282:G:O2'	23:BB:2283:C:OP2	2.15	0.61
23:BB:2308:G:H5'	23:BB:2309:A:H5''	1.82	0.61
23:BB:2471:A:O2'	23:BB:2472:G:O5'	2.19	0.61
23:BB:2839:G:H21	35:BN:92:GLY:HA3	1.64	0.61
23:BB:557:C:O2'	31:BJ:114:LEU:HD13	2.01	0.61
23:BB:620:G:N3	23:BB:620:G:H5'	2.15	0.61
23:BB:670:A:H3'	33:BL:47:ARG:NH1	2.10	0.61
25:BC:194:VAL:HG12	25:BC:195:GLY:N	2.16	0.61
26:BD:132:ALA:HA	26:BD:136:ASN:H	1.65	0.61
26:BD:21:SER:C	26:BD:23:PRO:HD3	2.20	0.61
33:BL:103:ILE:O	33:BL:104:GLN:HB3	2.00	0.61
33:BL:91:ASP:OD2	33:BL:123:ARG:HD2	2.01	0.61
34:BM:34:LYS:HA	34:BM:99:GLY:HA3	1.83	0.61
35:BN:98:LEU:O	35:BN:111:ALA:HA	2.01	0.61
44:BX:45:GLN:H	44:BX:45:GLN:HE21	1.49	0.61
1:CA:900:A:O2'	1:CA:901:A:H5'	2.00	0.61
20:CB:36:LYS:HE3	20:CB:36:LYS:N	2.16	0.61
9:CJ:64:GLN:HB3	13:CN:98:ALA:HB3	1.83	0.61
16:CQ:10:ARG:CZ	16:CQ:55:GLY:H	2.13	0.61
23:DB:2615:U:H1'	47:D0:3:GLN:HB3	1.83	0.61
22:DA:66:A:HO2'	22:DA:67:G:H8	1.49	0.61
23:DB:1019:U:H2'	23:DB:1020:A:C8	2.35	0.61
23:DB:1439:A:C4	23:DB:1552:A:N6	2.69	0.61
23:DB:291:G:O2'	23:DB:292:U:H5'	2.01	0.61
27:DE:99:LYS:HZ2	27:DE:100:MET:N	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:112:ARG:N	37:DP:112:ARG:HE	1.98	0.61
37:DP:23:ASP:HA	37:DP:49:ILE:CG2	2.30	0.61
39:DR:6:GLN:HE21	39:DR:6:GLN:N	1.98	0.61
1:AA:1137:C:O2'	1:AA:1138:G:H5''	2.00	0.61
1:AA:1280:A:O4'	9:AJ:43:PRO:HG3	2.01	0.61
1:AA:636:U:H2'	1:AA:637:C:H6	1.66	0.61
8:AI:51:LEU:HD11	8:AI:82:ILE:HD13	1.83	0.61
50:B3:41:ARG:HB3	50:B3:43:LEU:CD2	2.30	0.61
23:BB:1038:G:H2'	23:BB:1039:A:C8	2.36	0.61
23:BB:138:U:H2'	23:BB:140:C:H1'	1.83	0.61
23:BB:143:C:H2'	23:BB:144:A:C8	2.36	0.61
23:BB:2047:C:H2'	23:BB:2048:G:H8	1.66	0.61
23:BB:2318:G:O2'	23:BB:2319:G:N2	2.34	0.61
23:BB:224:U:O4	23:BB:420:C:H5'	2.00	0.61
23:BB:743:A:O3'	26:BD:136:ASN:HB2	2.00	0.61
23:BB:852:U:H2'	23:BB:853:C:C6	2.36	0.61
23:BB:969:G:H2'	23:BB:970:U:H6	1.64	0.61
26:BD:159:LYS:HB2	26:BD:159:LYS:HZ2	1.64	0.61
26:BD:165:MET:HE2	26:BD:166:GLY:H	1.66	0.61
27:BE:59:PRO:HD2	27:BE:60:TRP:NE1	2.16	0.61
28:BF:97:GLU:O	28:BF:100:GLU:HG3	2.00	0.61
33:BL:94:THR:O	33:BL:100:ILE:HG13	2.00	0.61
34:BM:17:ASN:C	34:BM:18:ARG:HD3	2.22	0.61
37:BP:52:ARG:HB3	37:BP:60:VAL:CG1	2.22	0.61
41:BT:45:ALA:O	41:BT:48:GLN:HB3	2.00	0.61
1:CA:56:U:H2'	1:CA:57:G:C8	2.36	0.61
3:CD:160:LEU:HD13	3:CD:161:ALA:N	2.16	0.61
4:CE:111:ARG:HG2	4:CE:111:ARG:HH11	1.66	0.61
1:CA:1319:A:OP1	18:CS:4:LEU:HD21	2.00	0.61
47:D0:36:LYS:CB	47:D0:41:HIS:HA	2.31	0.61
23:DB:1712:U:H3'	23:DB:1713:A:H5''	1.83	0.61
23:DB:1791:A:N6	23:DB:1828:G:H1'	2.16	0.61
23:DB:2455:G:H2'	23:DB:2456:C:C6	2.35	0.61
23:DB:2785:C:H2'	23:DB:2786:U:H6	1.65	0.61
23:DB:401:A:H2'	23:DB:402:A:C8	2.35	0.61
23:DB:898:C:C2'	23:DB:899:A:H5''	2.31	0.61
25:DC:12:ARG:HH11	25:DC:18:VAL:HB	1.66	0.61
27:DE:5:LEU:HA	27:DE:11:ALA:O	2.00	0.61
28:DF:133:GLU:HG2	28:DF:149:ARG:O	2.01	0.61
29:DG:36:LEU:CB	29:DG:40:VAL:HG21	2.30	0.61
38:DQ:108:LEU:HD12	38:DQ:111:LYS:HD2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1252:G:N3	38:DQ:32:ARG:HG3	2.16	0.61
41:DT:45:ALA:HA	41:DT:48:GLN:HG2	1.83	0.61
46:DZ:55:GLY:CA	46:DZ:59:ARG:HB2	2.26	0.61
1:AA:160:A:H2'	1:AA:161:A:O4'	2.00	0.60
1:AA:815:A:H4'	1:AA:817:C:C4	2.36	0.60
20:AB:67:LEU:HD21	20:AB:157:PRO:HG3	1.83	0.60
8:AI:87:MET:SD	8:AI:94:ARG:HG3	2.42	0.60
23:BB:1495:A:H2'	23:BB:1496:A:C8	2.35	0.60
23:BB:1821:A:H2'	23:BB:1822:C:C6	2.36	0.60
23:BB:2183:A:H2'	23:BB:2184:A:C8	2.36	0.60
23:BB:2895:G:H2'	23:BB:2896:C:H6	1.64	0.60
25:BC:253:GLY:C	25:BC:254:LYS:HD2	2.21	0.60
26:BD:128:ARG:HH22	26:BD:130:GLN:CD	2.05	0.60
26:BD:34:VAL:HG11	26:BD:71:ALA:HB2	1.83	0.60
52:BI:100:ILE:O	52:BI:139:VAL:HA	2.01	0.60
23:BB:1076:C:H4'	52:BI:94:LYS:HE3	1.81	0.60
31:BJ:100:VAL:HG22	31:BJ:101:ILE:N	2.15	0.60
33:BL:62:PRO:O	50:B3:24:LYS:HD3	2.01	0.60
36:BO:35:ILE:HD13	36:BO:35:ILE:N	2.16	0.60
38:BQ:33:VAL:HG13	38:BQ:34:ALA:N	2.09	0.60
39:BR:18:GLN:NE2	39:BR:99:THR:HB	2.15	0.60
23:BB:64:A:O2'	41:BT:69:ARG:HB3	2.01	0.60
1:CA:278:G:N2	1:CA:279:A:H62	1.97	0.60
20:CB:36:LYS:H	20:CB:36:LYS:HE3	1.66	0.60
12:CM:102:LYS:HZ2	12:CM:103:THR:HG23	1.66	0.60
22:DA:2:G:H2'	22:DA:3:C:C6	2.36	0.60
23:DB:1098:A:C3'	52:DI:3:LYS:C	2.69	0.60
23:DB:2339:C:H2'	23:DB:2340:A:C8	2.36	0.60
25:DC:174:ARG:HG3	25:DC:180:MET:HG3	1.83	0.60
25:DC:235:GLU:HG3	25:DC:237:ARG:HD3	1.82	0.60
26:DD:73:VAL:HB	26:DD:91:THR:HB	1.83	0.60
29:DG:36:LEU:H	29:DG:36:LEU:HD12	1.66	0.60
30:DH:1:MET:HG3	30:DH:21:VAL:CG2	2.31	0.60
52:DI:1:ALA:C	52:DI:2:LYS:HD2	2.21	0.60
31:DJ:49:ASP:HA	31:DJ:114:LEU:HD11	1.82	0.60
32:DK:70:ARG:HB3	32:DK:76:VAL:HA	1.82	0.60
35:DN:78:LYS:HG2	35:DN:82:GLU:HG3	1.83	0.60
38:DQ:88:GLU:CD	39:DR:53:PHE:HB2	2.21	0.60
42:DU:4:ILE:CG2	42:DU:27:VAL:HG22	2.31	0.60
1:AA:1053:G:H4'	1:AA:1054:C:H5'	1.83	0.60
1:AA:1142:G:H2'	1:AA:1143:G:O4'	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:122:ILE:O	3:AD:128:VAL:HG23	2.01	0.60
3:AD:19:PHE:HB2	3:AD:110:ARG:HH12	1.66	0.60
7:AH:64:TYR:HA	7:AH:70:VAL:HG23	1.82	0.60
10:AK:88:PRO:HA	10:AK:92:ARG:HE	1.65	0.60
23:BB:1474:U:H2'	23:BB:1475:G:H5'	1.84	0.60
23:BB:2196:C:O2'	23:BB:2197:U:H5'	2.01	0.60
23:BB:2298:A:N6	23:BB:2319:G:H22	1.99	0.60
23:BB:2760:C:O2'	23:BB:2761:A:H5'	2.01	0.60
23:BB:2785:C:H2'	23:BB:2786:U:H6	1.65	0.60
23:BB:445:C:O2'	23:BB:446:G:H5'	2.01	0.60
25:BC:92:LEU:O	25:BC:93:VAL:HB	2.02	0.60
31:BJ:84:ILE:HG22	31:BJ:85:LYS:N	2.16	0.60
33:BL:56:PRO:HA	33:BL:60:ARG:NH2	2.16	0.60
34:BM:109:PRO:HA	34:BM:112:LEU:HB2	1.83	0.60
35:BN:36:THR:O	35:BN:40:LYS:HB2	2.01	0.60
35:BN:9:GLN:NE2	35:BN:17:ARG:HH22	1.99	0.60
36:BO:87:ILE:CG1	36:BO:88:LYS:H	2.14	0.60
38:BQ:50:ARG:HA	38:BQ:53:LYS:HE3	1.83	0.60
23:BB:2270:A:H4'	43:BW:18:LYS:HB3	1.82	0.60
23:BB:2081:U:H5''	46:BZ:23:LYS:HD2	1.82	0.60
1:CA:1324:A:H5'	1:CA:1362:A:O2'	2.01	0.60
1:CA:57:G:H2'	1:CA:58:C:C6	2.35	0.60
1:CA:636:U:H2'	1:CA:637:C:H6	1.66	0.60
1:CA:642:A:H2'	1:CA:643:C:C6	2.35	0.60
1:CA:77:A:H2'	1:CA:78:A:H8	1.66	0.60
21:CU:40:PRO:O	21:CU:44:ARG:N	2.34	0.60
23:DB:1025:G:H1'	23:DB:1135:C:C5'	2.26	0.60
23:DB:2621:G:O2'	26:DD:164:GLN:HG3	2.00	0.60
23:DB:393:C:O2'	23:DB:394:C:H5'	2.01	0.60
23:DB:633:A:H2'	23:DB:634:C:H5'	1.82	0.60
23:DB:6:A:H2'	23:DB:7:G:H8	1.65	0.60
25:DC:117:SER:HB3	25:DC:127:ASN:ND2	2.17	0.60
23:DB:2680:U:H5'	26:DD:194:PRO:HA	1.83	0.60
27:DE:109:LEU:HG	27:DE:117:ARG:HD3	1.82	0.60
27:DE:14:VAL:HG12	27:DE:15:SER:N	2.16	0.60
30:DH:135:HIS:HB3	30:DH:138:VAL:HB	1.83	0.60
33:DL:78:ARG:CZ	33:DL:78:ARG:HB3	2.31	0.60
23:DB:2293:G:OP1	36:DO:13:ARG:NH2	2.34	0.60
37:DP:5:LYS:HA	37:DP:5:LYS:NZ	2.16	0.60
43:DW:39:GLN:HG2	43:DW:66:VAL:O	2.01	0.60
8:AI:51:LEU:HD23	8:AI:56:MET:CE	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AJ:15:HIS:HA	9:AJ:18:ILE:HG22	1.82	0.60
11:AL:66:ILE:HD13	11:AL:73:LEU:HD12	1.84	0.60
19:AT:4:LYS:HD2	19:AT:5:SER:N	2.16	0.60
48:B1:8:ILE:HD11	48:B1:25:ASN:HB3	1.82	0.60
23:BB:1857:G:H1'	23:BB:1885:A:N6	2.16	0.60
23:BB:1857:G:H2'	23:BB:1884:G:H22	1.66	0.60
23:BB:2785:C:H2'	23:BB:2786:U:C6	2.36	0.60
23:BB:28:A:N6	23:BB:512:G:O2'	2.33	0.60
23:BB:765:C:H2'	23:BB:766:U:H6	1.66	0.60
23:BB:903:C:H2'	23:BB:904:G:C8	2.37	0.60
26:BD:3:GLY:HA2	26:BD:49:GLN:OE1	2.01	0.60
28:BF:135:ILE:C	28:BF:137:PHE:H	2.04	0.60
52:BI:10:LEU:HD13	52:BI:12:VAL:HG13	1.82	0.60
36:BO:68:LYS:HD3	36:BO:68:LYS:N	2.16	0.60
39:BR:64:VAL:HB	39:BR:98:ILE:HD12	1.84	0.60
43:BW:45:HIS:HA	43:BW:75:ASN:CB	2.32	0.60
1:CA:1176:A:H2'	1:CA:1177:G:O4'	2.01	0.60
1:CA:26:A:H61	1:CA:558:G:H1'	1.65	0.60
4:CE:12:GLU:HG2	4:CE:38:VAL:HG22	1.83	0.60
7:CH:81:GLY:O	7:CH:82:LEU:HB2	2.02	0.60
11:CL:80:LEU:O	11:CL:97:VAL:HG22	2.00	0.60
23:DB:144:A:C2	41:DT:3:ARG:CZ	2.85	0.60
23:DB:2650:U:H2'	23:DB:2651:C:C6	2.36	0.60
23:DB:2789:C:H3'	23:DB:2893:A:H62	1.66	0.60
23:DB:2809:A:H2'	23:DB:2810:A:C8	2.35	0.60
23:DB:374:A:N6	23:DB:400:G:H1'	2.16	0.60
23:DB:533:G:H5'	38:DQ:23:TYR:CD1	2.37	0.60
25:DC:65:ASP:OD2	25:DC:101:ARG:HD3	2.01	0.60
27:DE:143:LEU:H	27:DE:143:LEU:HD22	1.66	0.60
28:DF:33:ILE:HD11	28:DF:99:PHE:HB2	1.83	0.60
32:DK:102:PRO:CB	32:DK:121:GLU:HG2	2.30	0.60
33:DL:96:LYS:HE3	33:DL:96:LYS:HA	1.81	0.60
35:DN:2:ARG:HG2	35:DN:2:ARG:NH1	2.15	0.60
39:DR:79:ARG:NH2	39:DR:87:GLN:HG3	2.15	0.60
41:DT:15:HIS:O	41:DT:16:VAL:HB	2.01	0.60
23:DB:2354:C:H4'	43:DW:30:VAL:HG13	1.82	0.60
46:DZ:64:PHE:O	46:DZ:67:PRO:HD2	2.01	0.60
1:AA:1526:G:O2'	1:AA:1527:U:H5'	2.01	0.60
1:AA:620:C:C6	3:AD:131:ILE:HD13	2.37	0.60
20:AB:113:LEU:O	20:AB:117:GLU:HG3	2.00	0.60
20:AB:116:LEU:HD11	20:AB:139:GLU:HB3	1.81	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AB:150:ILE:HG13	20:AB:153:MET:HE3	1.81	0.60
20:AB:204:ASP:CG	20:AB:205:ALA:H	2.03	0.60
12:AM:21:ILE:CG2	12:AM:64:VAL:HG11	2.30	0.60
13:AN:24:ALA:O	13:AN:27:LYS:HG3	2.01	0.60
18:AS:31:ARG:HG3	18:AS:56:HIS:NE2	2.15	0.60
19:AT:68:LYS:HG3	19:AT:69:ASN:N	2.10	0.60
23:BB:1439:A:C5	23:BB:1552:A:N6	2.70	0.60
23:BB:1810:A:H2'	23:BB:1811:G:O4'	2.00	0.60
23:BB:1939:U:O2	23:BB:1967:C:H4'	2.01	0.60
23:BB:2599:G:N7	25:BC:234:GLY:HA2	2.16	0.60
23:BB:2600:A:H62	25:BC:235:GLU:HB3	1.66	0.60
23:BB:2848:G:H2'	37:BP:96:LEU:HB2	1.83	0.60
23:BB:607:U:OP2	27:BE:98:LYS:HG2	2.02	0.60
23:BB:692:C:H5'	25:BC:42:ARG:CB	2.22	0.60
26:BD:31:ALA:HB3	26:BD:95:SER:HB3	1.84	0.60
27:BE:106:LYS:HZ2	27:BE:106:LYS:HB2	1.66	0.60
30:BH:116:ARG:HB3	30:BH:131:SER:OG	2.01	0.60
34:BM:3:GLN:H	34:BM:4:PRO:CD	1.95	0.60
35:BN:8:ARG:CZ	35:BN:46:ARG:HD2	2.32	0.60
42:BU:13:LEU:H	42:BU:68:ASN:ND2	2.00	0.60
1:CA:1129:C:H1'	1:CA:1146:A:H61	1.65	0.60
1:CA:815:A:H4'	1:CA:817:C:C4	2.36	0.60
1:CA:922:G:H2'	1:CA:923:A:H8	1.66	0.60
11:CL:29:LYS:O	11:CL:80:LEU:HD12	2.01	0.60
12:CM:15:VAL:HG22	12:CM:40:GLU:HB2	1.83	0.60
13:CN:79:SER:O	13:CN:83:VAL:HG23	2.02	0.60
14:CO:88:ARG:HH11	14:CO:88:ARG:HG3	1.66	0.60
15:CP:4:ILE:O	15:CP:71:VAL:HG11	2.00	0.60
48:D1:8:ILE:HD13	48:D1:9:LYS:H	1.64	0.60
51:D4:36:ARG:HG2	51:D4:36:ARG:HH11	1.66	0.60
23:DB:1810:A:H2'	23:DB:1811:G:O4'	2.00	0.60
23:DB:197:A:N6	23:DB:2430:A:H2'	2.17	0.60
23:DB:2888:C:H2'	23:DB:2889:C:C6	2.36	0.60
23:DB:352:A:H3'	23:DB:353:C:H6	1.66	0.60
23:DB:850:U:H2'	23:DB:851:C:C6	2.36	0.60
27:DE:176:ASP:HB3	27:DE:179:SER:OG	2.02	0.60
37:DP:36:LYS:HG2	37:DP:37:LYS:H	1.67	0.60
37:DP:70:GLU:HG2	37:DP:71:ARG:H	1.67	0.60
1:AA:57:G:H2'	1:AA:58:C:C6	2.36	0.60
1:AA:74:A:O2'	1:AA:75:G:H5''	2.02	0.60
1:AA:769:G:H4'	1:AA:1513:A:H4'	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:185:PRO:HB2	3:AD:190:LEU:HD11	1.81	0.60
7:AH:26:MET:HB2	7:AH:27:PRO:HD2	1.84	0.60
8:AI:110:VAL:HG12	8:AI:111:GLU:N	2.16	0.60
16:AQ:59:GLU:H	16:AQ:74:LEU:HD23	1.67	0.60
1:AA:1320:C:H5	18:AS:36:ARG:HG2	1.65	0.60
47:B0:36:LYS:HG3	47:B0:41:HIS:HA	1.83	0.60
23:BB:2331:G:H4'	43:BW:69:GLU:CG	2.31	0.60
25:BC:68:ARG:HD3	25:BC:103:ILE:CD1	2.31	0.60
33:BL:91:ASP:O	33:BL:92:LEU:CB	2.49	0.60
38:BQ:27:ARG:HG3	38:BQ:32:ARG:NH2	2.16	0.60
38:BQ:92:LYS:O	38:BQ:94:LEU:N	2.34	0.60
38:BQ:88:GLU:HA	39:BR:53:PHE:CE2	2.35	0.60
40:BS:32:ALA:C	40:BS:33:LEU:HD22	2.21	0.60
22:BA:98:G:H1	24:BV:14:LYS:CB	2.15	0.60
1:CA:403:C:O2'	1:CA:404:G:H5'	2.01	0.60
1:CA:60:A:O2'	1:CA:61:G:O5'	2.19	0.60
4:CE:91:SER:HA	4:CE:128:GLY:O	2.02	0.60
4:CE:36:THR:HG22	4:CE:37:VAL:N	2.14	0.60
5:CF:42:TRP:HE1	5:CF:61:LEU:HD23	1.67	0.60
13:CN:27:LYS:O	13:CN:31:SER:HB3	2.00	0.60
23:DB:2679:A:O2'	23:DB:2680:U:H5'	2.02	0.60
23:DB:2800:A:H2'	23:DB:2801:G:C1'	2.32	0.60
26:DD:107:VAL:HA	26:DD:205:PRO:O	2.01	0.60
23:DB:448:U:H3'	27:DE:79:ARG:NH2	2.15	0.60
52:DI:125:THR:O	52:DI:129:GLU:HG3	2.01	0.60
31:DJ:36:LEU:HA	31:DJ:51:GLY:O	2.02	0.60
44:DX:44:LYS:CG	44:DX:47:ARG:HB2	2.31	0.60
45:DY:6:ILE:CG1	45:DY:35:VAL:H	2.13	0.60
1:AA:859:G:H2'	1:AA:860:A:H8	1.65	0.60
2:AC:58:ARG:CG	2:AC:63:ILE:HG22	2.31	0.60
3:AD:150:LYS:HB2	3:AD:150:LYS:NZ	2.17	0.60
10:AK:52:ARG:HD2	10:AK:53:GLY:H	1.67	0.60
48:B1:5:ARG:HB3	48:B1:5:ARG:HH11	1.65	0.60
23:BB:1805:A:N3	25:BC:50:THR:HG23	2.16	0.60
23:BB:1878:G:H2'	23:BB:1879:C:C6	2.37	0.60
23:BB:2642:G:O2'	23:BB:2643:G:H5'	2.01	0.60
23:BB:2679:A:O2'	23:BB:2680:U:H5'	2.02	0.60
23:BB:2698:U:H2'	23:BB:2699:C:C6	2.37	0.60
23:BB:2769:U:H2'	23:BB:2770:G:C8	2.36	0.60
23:BB:328:U:H4'	42:BU:65:GLN:NE2	2.16	0.60
23:BB:926:G:H2'	23:BB:927:A:H8	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:30:ALA:N	25:BC:31:PRO:HD3	2.09	0.60
32:BK:2:ILE:HG22	32:BK:67:LYS:NZ	2.16	0.60
33:BL:77:ILE:CG2	33:BL:78:ARG:HG3	2.32	0.60
39:BR:74:ILE:CA	39:BR:90:ARG:HE	2.10	0.60
42:BU:85:ARG:N	42:BU:85:ARG:HD3	2.17	0.60
24:BV:46:LYS:HA	24:BV:46:LYS:HE3	1.84	0.60
1:CA:994:A:C5	1:CA:1216:A:H4'	2.36	0.60
20:CB:16:GLY:HA2	20:CB:40:ILE:H	1.66	0.60
3:CD:171:GLU:HG3	3:CD:182:LYS:HD2	1.83	0.60
4:CE:113:VAL:HG23	4:CE:114:LEU:H	1.66	0.60
23:DB:2887:A:C8	47:D0:27:LEU:HD21	2.36	0.60
22:DA:35:C:O2'	22:DA:36:C:H5'	2.02	0.60
23:DB:1082:U:O4	23:DB:1086:A:C2	2.55	0.60
23:DB:1098:A:C2'	52:DI:3:LYS:C	2.70	0.60
23:DB:990:A:N6	23:DB:1186:G:H1'	2.17	0.60
23:DB:1203:U:H3'	23:DB:1204:A:C5'	2.31	0.60
23:DB:1439:A:C5	23:DB:1552:A:N6	2.70	0.60
23:DB:1791:A:C2	23:DB:1829:A:H4'	2.36	0.60
23:DB:2360:G:H1'	33:DL:61:LEU:HD11	1.82	0.60
26:DD:24:VAL:HG13	26:DD:193:VAL:HG21	1.84	0.60
27:DE:157:LEU:H	27:DE:157:LEU:HD23	1.66	0.60
31:DJ:25:LEU:CD1	31:DJ:63:ALA:H	2.14	0.60
32:DK:9:ASN:O	32:DK:10:VAL:HG13	2.01	0.60
33:DL:38:GLN:NE2	33:DL:39:LYS:HE2	2.16	0.60
35:DN:56:LYS:CG	35:DN:57:THR:H	2.14	0.60
37:DP:48:ALA:O	37:DP:49:ILE:HG22	2.01	0.60
41:DT:88:LYS:HG3	41:DT:89:GLU:OE1	2.02	0.60
42:DU:78:LYS:HD2	42:DU:96:LYS:HG3	1.83	0.60
23:DB:2090:A:O2'	46:DZ:49:ARG:HD2	2.02	0.60
1:AA:812:G:HO2'	1:AA:813:U:H6	1.49	0.60
1:AA:812:G:O2'	1:AA:813:U:H6	1.85	0.60
5:AF:46:GLN:HA	5:AF:56:LYS:HG3	1.81	0.60
5:AF:61:LEU:HD12	5:AF:63:ASN:OD1	2.02	0.60
13:AN:15:LEU:HD12	13:AN:16:ALA:H	1.66	0.60
16:AQ:60:ILE:N	16:AQ:60:ILE:HD13	2.16	0.60
23:BB:176:A:O2'	23:BB:177:G:H5'	2.01	0.60
23:BB:1791:A:N6	23:BB:1828:G:H1'	2.16	0.60
23:BB:2650:U:H2'	23:BB:2651:C:C6	2.35	0.60
23:BB:2671:G:H2'	23:BB:2672:U:C6	2.37	0.60
23:BB:26:G:H1'	23:BB:514:A:N6	2.17	0.60
23:BB:418:C:H2'	23:BB:419:U:C6	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:433:C:H2'	23:BB:434:U:C6	2.37	0.60
23:BB:836:G:H2'	23:BB:837:C:C6	2.37	0.60
23:BB:910:A:N7	34:BM:16:ARG:HB2	2.17	0.60
25:BC:58:LYS:O	25:BC:59:GLN:HG3	2.02	0.60
28:BF:174:PHE:N	28:BF:175:PRO:CD	2.63	0.60
33:BL:60:ARG:HD2	50:B3:7:ARG:HE	1.66	0.60
37:BP:50:ARG:CB	37:BP:50:ARG:HH11	2.09	0.60
23:BB:997:G:H5'	38:BQ:91:ARG:HG3	1.83	0.60
40:BS:29:VAL:HG22	40:BS:71:VAL:HG23	1.83	0.60
44:BX:44:LYS:O	44:BX:48:ARG:HD2	2.02	0.60
1:CA:1256:A:N1	1:CA:1278:G:H1'	2.16	0.60
1:CA:590:U:H2'	1:CA:591:U:H6	1.67	0.60
2:CC:11:LEU:HD22	2:CC:17:TRP:NE1	2.17	0.60
2:CC:56:ILE:N	2:CC:56:ILE:HD12	2.16	0.60
4:CE:142:GLY:HA2	4:CE:145:ASN:ND2	2.16	0.60
5:CF:36:ILE:HG12	5:CF:64:VAL:HG22	1.83	0.60
1:CA:1124:G:H5'	9:CJ:37:ARG:NH2	2.17	0.60
13:CN:30:ILE:CG2	13:CN:41:TRP:HB3	2.31	0.60
23:DB:1857:G:H2'	23:DB:1884:G:H22	1.66	0.60
23:DB:1878:G:H2'	23:DB:1879:C:C6	2.37	0.60
23:DB:1915:U:H2'	23:DB:1916:A:O4'	2.02	0.60
23:DB:1939:U:O2	23:DB:1967:C:H4'	2.01	0.60
25:DC:92:LEU:O	25:DC:93:VAL:HB	2.01	0.60
26:DD:129:THR:HA	26:DD:140:HIS:CE1	2.36	0.60
33:DL:91:ASP:O	33:DL:93:ASN:N	2.33	0.60
34:DM:29:GLY:N	34:DM:102:LEU:HD12	2.12	0.60
35:DN:54:LEU:O	35:DN:54:LEU:HD23	2.02	0.60
41:DT:61:LEU:HB2	41:DT:82:LYS:HB3	1.83	0.60
44:DX:30:MET:H	44:DX:30:MET:HE2	1.67	0.60
1:AA:51:A:O2'	1:AA:52:C:OP2	2.18	0.60
1:AA:26:A:H61	1:AA:558:G:H1'	1.65	0.60
1:AA:56:U:H2'	1:AA:57:G:C8	2.36	0.60
1:AA:590:U:H2'	1:AA:591:U:H6	1.66	0.60
20:AB:46:VAL:CG1	20:AB:47:PRO:HD3	2.25	0.60
4:AE:109:ALA:HB3	4:AE:135:VAL:HG23	1.82	0.60
4:AE:85:LYS:HE3	4:AE:94:PHE:HB2	1.84	0.60
8:AI:49:GLN:HE21	8:AI:79:ARG:HD2	1.67	0.60
17:AR:46:THR:HG23	17:AR:51:GLN:HB2	1.81	0.60
23:BB:1729:U:C5	23:BB:1730:C:H1'	2.37	0.60
23:BB:27:G:H1'	23:BB:513:A:N6	2.17	0.60
23:BB:575:A:O2'	23:BB:576:U:H5'	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2050:C:H4'	26:BD:143:PRO:HG2	1.84	0.60
26:BD:153:GLY:O	26:BD:154:LYS:HB2	2.01	0.60
30:BH:6:LEU:HA	30:BH:15:LEU:HA	1.83	0.60
33:BL:126:ARG:O	33:BL:127:VAL:HG13	2.02	0.60
36:BO:16:ARG:O	36:BO:20:GLU:HB2	2.01	0.60
40:BS:24:ILE:HD13	40:BS:36:LEU:HD21	1.83	0.60
42:BU:12:VAL:HA	42:BU:68:ASN:ND2	2.12	0.60
23:BB:2365:G:H4'	43:BW:65:LYS:HZ2	1.67	0.60
1:CA:1118:U:H2'	1:CA:1119:C:H6	1.67	0.60
1:CA:135:C:O2	15:CP:1:MET:HB2	2.02	0.60
2:CC:110:LEU:HD13	2:CC:143:LEU:HD23	1.83	0.60
4:CE:158:LYS:CB	7:CH:63:LYS:HD3	2.30	0.60
13:CN:30:ILE:HB	13:CN:44:VAL:HB	1.83	0.60
18:CS:27:LYS:HB3	18:CS:27:LYS:NZ	2.17	0.60
1:CA:1320:C:OP2	18:CS:2:ARG:HB3	2.02	0.60
23:DB:1164:C:H2'	23:DB:1165:A:H8	1.65	0.60
23:DB:2868:A:H2'	23:DB:2869:G:C8	2.37	0.60
23:DB:717:C:C3'	23:DB:718:A:H5''	2.30	0.60
25:DC:257:ARG:HH11	25:DC:257:ARG:HG2	1.66	0.60
52:DI:63:ASP:O	52:DI:64:ARG:HB2	2.00	0.60
34:DM:41:LEU:HB3	34:DM:93:VAL:HB	1.83	0.60
37:DP:106:ALA:O	37:DP:108:ARG:HD2	2.01	0.60
38:DQ:98:ALA:HA	38:DQ:105:PHE:CD1	2.36	0.60
41:DT:24:MET:HE1	41:DT:30:ILE:HG13	1.83	0.60
46:DZ:30:HIS:O	46:DZ:48:GLN:HA	2.01	0.60
46:DZ:1:MET:HA	46:DZ:9:TYR:CD1	2.37	0.60
1:AA:34:C:H2'	1:AA:35:G:H8	1.67	0.60
1:AA:865:A:H2'	1:AA:866:C:H6	1.67	0.60
1:AA:987:G:H2'	1:AA:988:G:C8	2.35	0.60
20:AB:162:VAL:HG13	20:AB:184:ALA:HB2	1.83	0.60
2:AC:45:GLU:C	2:AC:46:LEU:HD12	2.22	0.60
2:AC:5:HIS:CD2	2:AC:8:GLY:H	2.19	0.60
5:AF:37:HIS:CE1	5:AF:65:GLU:HB2	2.36	0.60
8:AI:18:VAL:HG22	8:AI:64:ILE:HG23	1.84	0.60
12:AM:103:THR:HG22	12:AM:104:ASN:H	1.67	0.60
51:B4:24:ARG:HH12	51:B4:36:ARG:HB2	1.66	0.60
23:BB:1082:U:O4	23:BB:1086:A:C2	2.55	0.60
23:BB:1655:A:C2	23:BB:2049:G:H5''	2.36	0.60
23:BB:2674:G:H4'	32:BK:30:ARG:CG	2.31	0.60
25:BC:32:LEU:HD13	25:BC:61:TYR:HE1	1.67	0.60
26:BD:5:VAL:HG11	26:BD:28:GLU:CA	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:BJ:90:GLU:O	31:BJ:93:ILE:HG22	2.02	0.60
33:BL:39:LYS:HB2	33:BL:46:VAL:HA	1.83	0.60
33:BL:83:ALA:N	33:BL:86:GLU:HG3	2.15	0.60
1:CA:1513:A:H2'	1:CA:1514:G:C8	2.37	0.60
1:CA:160:A:H2'	1:CA:161:A:O4'	2.00	0.60
20:CB:187:ASP:HB3	20:CB:201:GLY:O	2.02	0.60
4:CE:19:ARG:HG2	4:CE:20:VAL:N	2.15	0.60
7:CH:75:GLN:NE2	7:CH:76:ARG:H	1.99	0.60
18:CS:18:VAL:HG22	18:CS:42:ASN:ND2	2.17	0.60
23:DB:138:U:H2'	23:DB:140:C:O4'	2.01	0.60
23:DB:2386:A:H2'	23:DB:2387:U:C6	2.37	0.60
23:DB:2698:U:H2'	23:DB:2699:C:C6	2.36	0.60
26:DD:37:VAL:HG13	26:DD:42:ASN:CB	2.31	0.60
27:DE:105:LEU:HD22	27:DE:105:LEU:H	1.67	0.60
27:DE:76:PRO:HA	27:DE:82:GLY:O	2.02	0.60
29:DG:138:GLN:NE2	29:DG:142:GLN:HG3	2.17	0.60
31:DJ:85:LYS:HA	31:DJ:85:LYS:HE2	1.83	0.60
37:DP:47:ILE:CG2	37:DP:63:ILE:HA	2.28	0.60
37:DP:47:ILE:CG1	37:DP:63:ILE:HG23	2.26	0.60
23:DB:992:C:H4'	38:DQ:46:TYR:OH	2.01	0.60
24:DV:77:VAL:HG13	24:DV:89:ILE:HD11	1.82	0.60
1:AA:1458:G:H5'	19:AT:26:MET:HB2	1.84	0.60
1:AA:238:A:C2'	1:AA:239:U:H5''	2.31	0.60
7:AH:93:LYS:H	7:AH:93:LYS:HZ2	1.49	0.60
10:AK:86:LYS:HD2	10:AK:114:PRO:HD3	1.84	0.60
11:AL:85:ARG:HA	11:AL:93:ARG:HA	1.84	0.60
49:B2:17:GLY:HA2	49:B2:21:ARG:NH1	2.17	0.60
49:B2:34:ARG:HE	49:B2:42:LEU:CD1	2.14	0.60
22:BA:39:A:O2'	22:BA:40:U:H5'	2.01	0.60
22:BA:75:G:H5''	24:BV:12:GLN:OE1	2.01	0.60
23:BB:1557:C:H3'	23:BB:1558:C:H5''	1.84	0.60
23:BB:2811:G:C5'	26:BD:62:LYS:HD2	2.31	0.60
23:BB:634:C:H2'	23:BB:635:C:C6	2.37	0.60
23:BB:651:G:H4'	50:B3:17:GLY:HA3	1.84	0.60
23:BB:864:G:O2'	23:BB:865:C:H5'	2.02	0.60
28:BF:64:PRO:HA	28:BF:88:VAL:CG2	2.32	0.60
30:BH:21:VAL:HG22	30:BH:22:LYS:N	2.13	0.60
33:BL:120:VAL:HG22	33:BL:122:VAL:HG13	1.83	0.60
23:BB:1454:C:C5'	35:BN:63:ARG:HH21	2.06	0.60
38:BQ:94:LEU:HD22	38:BQ:95:ALA:N	2.16	0.60
35:BN:107:ASN:OD1	40:BS:40:ASN:HB3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BT:24:MET:CE	41:BT:29:THR:H	2.15	0.60
42:BU:42:LYS:HD3	42:BU:60:LYS:HB3	1.83	0.60
24:BV:26:PHE:HE2	24:BV:44:HIS:HA	1.66	0.60
1:CA:1038:C:H2'	1:CA:1039:G:C8	2.36	0.60
1:CA:599:C:H5''	7:CH:86:LYS:O	2.02	0.60
7:CH:86:LYS:HB3	7:CH:90:GLU:O	2.02	0.60
9:CJ:22:THR:O	9:CJ:26:VAL:HG23	2.01	0.60
9:CJ:42:LEU:HB3	9:CJ:43:PRO:HD2	1.84	0.60
12:CM:52:ILE:HD12	12:CM:55:LEU:HD12	1.83	0.60
13:CN:62:ARG:HH21	13:CN:69:PRO:HG3	1.65	0.60
1:CA:254:G:OP1	16:CQ:68:LYS:O	2.20	0.60
23:DB:1429:G:H2'	23:DB:1430:G:C8	2.37	0.60
23:DB:1484:U:H2'	23:DB:1485:U:C6	2.37	0.60
23:DB:2241:A:H2'	23:DB:2242:G:C8	2.36	0.60
23:DB:2671:G:H2'	23:DB:2672:U:C6	2.36	0.60
52:DI:11:GLN:HA	52:DI:55:PRO:HA	1.83	0.60
33:DL:33:ARG:HB3	39:DR:85:LYS:HZ1	1.65	0.60
23:DB:2010:G:H5''	40:DS:42:LYS:CB	2.31	0.60
44:DX:58:ASN:ND2	44:DX:63:ALA:HB3	2.16	0.60
45:DY:18:LYS:O	45:DY:22:THR:HG23	2.02	0.60
1:AA:1035:A:H2'	1:AA:1036:A:C8	2.36	0.59
1:AA:320:A:H2'	1:AA:321:A:C8	2.37	0.59
20:AB:118:THR:O	20:AB:121:GLN:HB3	2.02	0.59
3:AD:194:ILE:HD13	3:AD:195:ASN:H	1.67	0.59
3:AD:58:GLN:HA	3:AD:61:ARG:HB2	1.84	0.59
16:AQ:24:ILE:HB	16:AQ:41:THR:HB	1.82	0.59
23:BB:1229:C:H2'	23:BB:1230:A:H8	1.65	0.59
23:BB:1295:C:H2'	23:BB:1296:G:C8	2.36	0.59
23:BB:401:A:H2'	23:BB:402:A:C8	2.37	0.59
23:BB:624:C:OP1	50:B3:61:LEU:HB3	2.01	0.59
23:BB:956:G:N2	23:BB:959:A:H3'	2.16	0.59
25:BC:143:VAL:O	25:BC:152:GLN:HB2	2.01	0.59
25:BC:61:TYR:CD2	25:BC:84:PRO:HD2	2.37	0.59
26:BD:35:THR:HG22	26:BD:46:ARG:NH2	2.16	0.59
26:BD:49:GLN:O	26:BD:78:GLY:HA2	2.00	0.59
27:BE:145:ASP:C	27:BE:147:LEU:H	2.04	0.59
33:BL:90:VAL:HG11	33:BL:121:THR:N	2.11	0.59
41:BT:64:LYS:NZ	41:BT:79:ASP:HB2	2.17	0.59
1:CA:1021:A:H2'	1:CA:1022:A:O4'	2.02	0.59
1:CA:1355:G:H2'	1:CA:1356:G:H8	1.67	0.59
1:CA:320:A:H2'	1:CA:321:A:C8	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:742:G:O2'	1:CA:743:A:H5'	2.02	0.59
1:CA:77:A:H2'	1:CA:78:A:C8	2.36	0.59
1:CA:22:G:H4'	1:CA:885:G:C8	2.37	0.59
3:CD:101:VAL:HG13	3:CD:106:PHE:CD2	2.34	0.59
4:CE:152:VAL:HG21	7:CH:98:LEU:HB3	1.82	0.59
4:CE:28:ARG:HH22	4:CE:30:PHE:HB3	1.67	0.59
7:CH:76:ARG:HE	7:CH:125:ILE:HG23	1.67	0.59
7:CH:31:LEU:O	7:CH:35:ILE:HG13	2.01	0.59
13:CN:26:LEU:HD12	13:CN:29:ILE:HD12	1.83	0.59
19:CT:66:ILE:O	19:CT:70:LYS:HD3	2.02	0.59
49:D2:35:ARG:NH2	49:D2:44:VAL:HG22	2.17	0.59
23:DB:1387:A:H5'	23:DB:1469:A:H1'	1.84	0.59
23:DB:1474:U:H2'	23:DB:1475:G:H5'	1.83	0.59
23:DB:2617:U:H2'	23:DB:2618:G:H5'	1.84	0.59
23:DB:2769:U:H2'	23:DB:2770:G:C8	2.36	0.59
23:DB:322:A:H1'	23:DB:339:U:O2	2.01	0.59
23:DB:493:G:H2'	23:DB:494:G:O4'	2.01	0.59
23:DB:858:G:H21	23:DB:2268:A:H3'	1.67	0.59
25:DC:143:VAL:HG11	25:DC:173:LEU:HD11	1.84	0.59
26:DD:125:TRP:HB2	26:DD:160:LYS:HD3	1.84	0.59
27:DE:187:VAL:HG23	27:DE:188:MET:N	2.17	0.59
30:DH:124:THR:HG23	30:DH:128:HIS:NE2	2.17	0.59
32:DK:11:ALA:HB1	32:DK:100:PHE:O	2.01	0.59
1:AA:1137:C:H1'	1:AA:1138:G:N2	2.16	0.59
1:AA:1103:C:H5''	20:AB:96:LEU:HD12	1.82	0.59
8:AI:56:MET:HA	8:AI:59:LYS:HZ2	1.67	0.59
23:BB:1285:A:H2'	23:BB:1286:A:H5''	1.83	0.59
23:BB:1550:C:H2'	23:BB:1551:A:H8	1.67	0.59
23:BB:2226:C:H2'	23:BB:2227:A:O4'	2.02	0.59
23:BB:2859:G:H2'	23:BB:2860:A:C8	2.36	0.59
23:BB:28:A:N6	23:BB:512:G:H1'	2.17	0.59
25:BC:128:THR:HG22	25:BC:188:ARG:HB2	1.84	0.59
25:BC:77:VAL:HG23	25:BC:115:ILE:CD1	2.31	0.59
28:BF:132:ARG:CZ	28:BF:147:ARG:HE	2.15	0.59
30:BH:125:THR:CB	30:BH:146:VAL:HB	2.31	0.59
31:BJ:28:LEU:HG	31:BJ:32:LEU:HD13	1.84	0.59
34:BM:102:LEU:HB3	34:BM:104:GLU:OE2	2.02	0.59
35:BN:96:ARG:NH2	35:BN:118:ARG:HB3	2.17	0.59
37:BP:30:TRP:HB2	37:BP:82:SER:OG	2.01	0.59
24:BV:14:LYS:NZ	24:BV:18:ARG:HD2	2.18	0.59
46:BZ:5:ILE:HD13	46:BZ:47:LYS:HB3	1.82	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1095:U:H2'	1:CA:1096:C:C6	2.37	0.59
6:CG:2:ARG:HH11	6:CG:2:ARG:HB3	1.66	0.59
13:CN:51:PRO:HB2	13:CN:54:SER:HB3	1.84	0.59
48:D1:13:SER:HB3	48:D1:49:LYS:NZ	2.17	0.59
23:DB:1729:U:C5	23:DB:1730:C:H1'	2.37	0.59
23:DB:1799:G:OP1	25:DC:255:LYS:HG2	2.02	0.59
23:DB:1874:C:H2'	23:DB:1875:G:O4'	2.01	0.59
23:DB:575:A:O2'	23:DB:576:U:H5'	2.02	0.59
23:DB:969:G:H2'	23:DB:970:U:H6	1.66	0.59
29:DG:19:ASN:HB3	29:DG:22:VAL:HB	1.84	0.59
30:DH:2:GLN:HA	30:DH:21:VAL:HG13	1.82	0.59
35:DN:96:ARG:O	35:DN:113:ILE:HA	2.01	0.59
36:DO:50:ALA:HB1	36:DO:78:VAL:HG13	1.82	0.59
37:DP:18:SER:HB2	37:DP:87:ARG:CZ	2.31	0.59
37:DP:25:VAL:HG22	37:DP:89:GLY:N	2.15	0.59
40:DS:21:ALA:HB1	40:DS:74:ILE:CD1	2.32	0.59
40:DS:85:ILE:HD11	40:DS:93:ALA:HB1	1.82	0.59
43:DW:48:ALA:HA	43:DW:54:ARG:H	1.67	0.59
43:DW:46:ALA:HB2	43:DW:75:ASN:O	2.02	0.59
8:AI:53:LEU:O	8:AI:53:LEU:HD13	2.02	0.59
23:BB:2539:C:OP1	51:B4:4:ARG:HD2	2.02	0.59
51:B4:4:ARG:HG2	51:B4:4:ARG:O	2.02	0.59
23:BB:143:C:H3'	23:BB:144:A:C8	2.37	0.59
23:BB:2273:A:H2'	23:BB:2274:A:C8	2.37	0.59
23:BB:2339:C:H2'	23:BB:2340:A:C8	2.37	0.59
23:BB:660:C:H2'	23:BB:661:A:H8	1.68	0.59
23:BB:672:C:O2'	23:BB:673:C:H5'	2.03	0.59
27:BE:149:ILE:HG13	27:BE:186:VAL:HG13	1.84	0.59
28:BF:132:ARG:HD3	28:BF:133:GLU:HB2	1.84	0.59
29:BG:157:LYS:HG3	29:BG:159:LYS:NZ	2.16	0.59
31:BJ:43:GLU:HB3	38:BQ:99:VAL:HG12	1.84	0.59
42:BU:48:VAL:HG12	42:BU:50:ALA:HB3	1.84	0.59
1:CA:202:G:H2'	1:CA:203:G:C8	2.37	0.59
20:CB:96:LEU:O	20:CB:99:MET:HG3	2.02	0.59
4:CE:96:GLN:HG3	4:CE:97:PRO:HD2	1.84	0.59
6:CG:12:LEU:HD13	6:CG:13:PRO:CD	2.31	0.59
10:CK:53:GLY:O	10:CK:56:LYS:HB3	2.02	0.59
12:CM:10:ASP:HA	12:CM:44:ILE:HD11	1.83	0.59
23:DB:585:G:H2'	23:DB:1251:C:H42	1.66	0.59
23:DB:136:G:N2	41:DT:3:ARG:HG2	2.17	0.59
23:DB:1695:G:O2'	25:DC:15:VAL:HG23	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1814:G:H4'	25:DC:51:ARG:HG2	1.83	0.59
23:DB:2107:G:H2'	23:DB:2108:A:H8	1.67	0.59
23:DB:2785:C:H2'	23:DB:2786:U:C6	2.37	0.59
23:DB:2789:C:H2'	23:DB:2893:A:N7	2.17	0.59
23:DB:2900:A:H2'	23:DB:2901:C:H6	1.66	0.59
25:DC:10:PRO:HD2	25:DC:202:ARG:NH2	2.16	0.59
27:DE:109:LEU:HD21	27:DE:113:VAL:O	2.02	0.59
29:DG:17:LYS:HE2	29:DG:24:THR:N	2.18	0.59
30:DH:115:VAL:HB	30:DH:132:PHE:CD1	2.37	0.59
30:DH:40:THR:HB	46:DZ:58:ASP:OD2	2.03	0.59
32:DK:34:GLY:O	32:DK:37:ASP:HB2	2.03	0.59
36:DO:100:HIS:HA	36:DO:104:GLN:HE21	1.67	0.59
40:DS:8:ARG:O	40:DS:9:HIS:HB2	2.02	0.59
1:AA:1060:U:H2'	1:AA:1061:G:C8	2.37	0.59
1:AA:1452:C:H4'	1:AA:1453:G:H5''	1.84	0.59
1:AA:532:A:H62	2:AC:191:THR:CG2	2.16	0.59
3:AD:94:GLU:HA	3:AD:103:ARG:HH22	1.66	0.59
9:AJ:26:VAL:O	9:AJ:30:LYS:HB3	2.02	0.59
1:AA:1060:U:C5'	9:AJ:53:ILE:HG22	2.32	0.59
19:AT:43:LYS:HG2	19:AT:44:ALA:N	2.17	0.59
23:BB:1054:A:H2'	23:BB:1055:G:H8	1.66	0.59
23:BB:182:A:H2'	23:BB:183:C:C6	2.37	0.59
23:BB:2266:A:H4'	23:BB:2267:A:C8	2.37	0.59
23:BB:2852:G:H2'	23:BB:2853:C:H6	1.67	0.59
23:BB:2868:A:H2'	23:BB:2869:G:C8	2.37	0.59
23:BB:2899:A:H2'	23:BB:2900:A:C8	2.37	0.59
23:BB:354:A:H2'	23:BB:355:U:H5''	1.83	0.59
23:BB:503:A:C4'	23:BB:504:A:H5'	2.31	0.59
23:BB:634:C:H2'	23:BB:635:C:H6	1.67	0.59
23:BB:784:G:O2'	23:BB:785:G:H5''	2.00	0.59
29:BG:85:LYS:HB2	29:BG:164:ALA:HB2	1.83	0.59
32:BK:23:LYS:HG3	32:BK:24:VAL:N	2.17	0.59
33:BL:117:THR:HG21	33:BL:120:VAL:HB	1.83	0.59
36:BO:15:ARG:HH12	36:BO:17:LYS:HD3	1.67	0.59
39:BR:90:ARG:NH2	39:BR:92:TRP:HB3	2.17	0.59
44:BX:49:ASP:HB3	44:BX:52:ARG:NH2	2.17	0.59
1:CA:1329:A:H5''	12:CM:25:GLY:N	2.16	0.59
1:CA:203:G:N2	1:CA:205:A:H61	2.01	0.59
1:CA:214:C:H2'	1:CA:215:C:C6	2.38	0.59
1:CA:621:A:H2'	1:CA:622:A:C8	2.37	0.59
2:CC:129:PHE:HA	2:CC:132:ALA:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:115:GLN:HE22	3:CD:153:ARG:HH22	1.50	0.59
9:CJ:36:VAL:HG13	9:CJ:76:ILE:HD11	1.83	0.59
23:DB:1285:A:H2'	23:DB:1286:A:H5''	1.83	0.59
23:DB:1295:C:H2'	23:DB:1296:G:C8	2.36	0.59
23:DB:1469:A:H2'	23:DB:1470:A:H8	1.63	0.59
23:DB:1676:A:H2'	23:DB:1677:A:O4'	2.02	0.59
23:DB:235:U:H2'	23:DB:236:C:H6	1.67	0.59
23:DB:533:G:H21	38:DQ:40:LYS:HD2	1.67	0.59
23:DB:544:C:H2'	23:DB:545:U:C5	2.37	0.59
23:DB:899:A:H2'	23:DB:900:A:O4'	2.02	0.59
25:DC:225:ASN:O	25:DC:227:VAL:N	2.35	0.59
27:DE:46:GLN:HE22	27:DE:86:ALA:HB3	1.68	0.59
12:CM:78:ARG:HG3	28:DF:111:ARG:HG2	1.83	0.59
30:DH:1:MET:O	30:DH:21:VAL:HG22	2.01	0.59
32:DK:41:ILE:HG13	32:DK:42:THR:N	2.17	0.59
32:DK:8:LEU:HD12	32:DK:8:LEU:N	2.17	0.59
34:DM:71:LYS:HB2	34:DM:72:PRO:CD	2.32	0.59
35:DN:32:GLU:HG3	35:DN:33:ILE:N	2.17	0.59
23:DB:534:U:H4'	38:DQ:45:ALA:HB2	1.83	0.59
39:DR:92:TRP:O	39:DR:93:PHE:HB2	2.01	0.59
1:AA:1055:A:H4'	2:AC:160:GLU:CG	2.32	0.59
1:AA:1101:A:O2'	1:AA:1102:A:OP2	2.21	0.59
1:AA:131:A:H2'	1:AA:132:C:H6	1.68	0.59
20:AB:18:GLN:O	20:AB:37:VAL:HG23	2.02	0.59
3:AD:7:LYS:HA	3:AD:7:LYS:HE3	1.85	0.59
7:AH:9:MET:HE1	7:AH:32:LYS:O	2.03	0.59
13:AN:20:PHE:HA	13:AN:24:ALA:HB2	1.84	0.59
15:AP:4:ILE:O	15:AP:71:VAL:HG11	2.02	0.59
21:AU:36:PHE:CB	21:AU:40:PRO:HD3	2.31	0.59
47:B0:39:ARG:HG2	47:B0:39:ARG:O	2.03	0.59
23:BB:1818:U:OP1	25:BC:155:ARG:HA	2.03	0.59
23:BB:2800:A:H2'	23:BB:2801:G:C1'	2.32	0.59
23:BB:478:A:H5''	23:BB:479:A:OP2	2.02	0.59
23:BB:718:A:H3'	23:BB:719:C:H6	1.68	0.59
25:BC:242:HIS:N	25:BC:243:PRO:CD	2.66	0.59
26:BD:38:LYS:H	26:BD:38:LYS:CE	2.15	0.59
29:BG:5:LYS:HB3	29:BG:68:ARG:NH2	2.17	0.59
31:BJ:136:GLN:N	31:BJ:137:PRO:CD	2.66	0.59
23:BB:1132:U:H5''	31:BJ:85:LYS:HZ3	1.67	0.59
34:BM:33:LEU:HD12	34:BM:101:VAL:HG11	1.83	0.59
23:BB:2821:A:OP2	35:BN:3:HIS:NE2	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BO:87:ILE:HG13	36:BO:88:LYS:N	2.16	0.59
44:BX:23:ARG:HG3	44:BX:28:LEU:HG	1.84	0.59
1:CA:1082:A:O2'	1:CA:1083:U:H5'	2.02	0.59
1:CA:1137:C:O2'	1:CA:1138:G:H5''	2.03	0.59
7:CH:76:ARG:HA	7:CH:126:CYS:HB3	1.85	0.59
16:CQ:14:ASP:OD2	16:CQ:53:GLY:HA2	2.01	0.59
16:CQ:10:ARG:NH2	16:CQ:55:GLY:H	2.00	0.59
23:DB:1097:U:H2'	23:DB:1098:A:O4'	2.03	0.59
23:DB:1945:G:H2'	23:DB:1946:U:C6	2.38	0.59
23:DB:2329:U:H2'	23:DB:2330:G:C8	2.37	0.59
23:DB:2642:G:O2'	23:DB:2643:G:H5'	2.01	0.59
25:DC:136:VAL:CA	25:DC:165:ALA:HA	2.33	0.59
25:DC:138:SER:HA	25:DC:162:GLN:HG3	1.85	0.59
26:DD:3:GLY:C	26:DD:4:LEU:HD13	2.23	0.59
28:DF:53:ALA:HB1	28:DF:64:PRO:HG2	1.84	0.59
29:DG:148:ARG:HA	29:DG:161:VAL:HG13	1.85	0.59
23:DB:1099:G:C8	52:DI:3:LYS:CA	2.75	0.59
38:DQ:54:ARG:O	38:DQ:58:GLN:HG2	2.02	0.59
40:DS:25:ARG:HD2	40:DS:26:GLY:H	1.65	0.59
43:DW:44:PHE:HB3	43:DW:77:LYS:CB	2.32	0.59
45:DY:6:ILE:HG13	45:DY:35:VAL:H	1.66	0.59
1:AA:1042:A:H2'	1:AA:1043:G:O4'	2.02	0.59
1:AA:214:C:H2'	1:AA:215:C:C6	2.38	0.59
2:AC:13:ILE:HG22	2:AC:14:VAL:HG13	1.83	0.59
4:AE:37:VAL:HG11	4:AE:113:VAL:HG12	1.83	0.59
22:BA:21:G:O2'	22:BA:22:U:H5'	2.02	0.59
22:BA:29:A:H3'	22:BA:30:C:H6	1.67	0.59
22:BA:24:G:N7	22:BA:56:G:H2'	2.17	0.59
23:BB:1412:U:H2'	23:BB:1413:A:C8	2.38	0.59
23:BB:1444:G:H2'	23:BB:1445:G:H8	1.67	0.59
23:BB:1676:A:H2'	23:BB:1677:A:O4'	2.02	0.59
23:BB:2077:A:H2'	23:BB:2078:C:C6	2.36	0.59
23:BB:2893:A:H5''	23:BB:2894:G:H5'	1.85	0.59
23:BB:359:G:O2'	23:BB:360:U:H5'	2.02	0.59
23:BB:79:C:O2'	23:BB:346:A:H1'	2.02	0.59
26:BD:4:LEU:O	26:BD:5:VAL:HG13	2.02	0.59
28:BF:115:GLY:HA3	28:BF:177:ARG:HB2	1.84	0.59
29:BG:131:VAL:HG13	29:BG:133:LYS:HE2	1.85	0.59
31:BJ:55:ILE:HA	31:BJ:122:LEU:O	2.02	0.59
33:BL:13:LYS:O	33:BL:14:LYS:HB2	2.01	0.59
35:BN:45:ARG:HB2	35:BN:49:GLU:OE2	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:77:SER:H	37:BP:78:PRO:HD2	1.68	0.59
23:BB:1162:G:H4'	39:BR:23:GLU:HG2	1.83	0.59
38:BQ:88:GLU:HA	39:BR:53:PHE:CZ	2.36	0.59
40:BS:52:GLU:HA	40:BS:55:ILE:HG22	1.85	0.59
1:CA:1030:U:H5''	1:CA:1031:C:H5	1.68	0.59
1:CA:1107:C:H4'	2:CC:172:VAL:HG23	1.83	0.59
1:CA:1251:A:O2'	1:CA:1252:A:H5'	2.02	0.59
1:CA:1308:U:H2'	1:CA:1309:G:C8	2.38	0.59
20:CB:116:LEU:O	20:CB:119:GLN:HB2	2.02	0.59
10:CK:53:GLY:HA2	10:CK:56:LYS:HE3	1.84	0.59
10:CK:82:GLU:HB3	10:CK:108:ASN:ND2	2.16	0.59
13:CN:96:LYS:HG2	13:CN:97:LYS:N	2.14	0.59
15:CP:4:ILE:HD13	15:CP:57:ILE:HG12	1.85	0.59
15:CP:57:ILE:C	15:CP:57:ILE:HD13	2.22	0.59
16:CQ:16:MET:CB	16:CQ:19:SER:HB2	2.33	0.59
17:CR:55:ALA:O	17:CR:59:LYS:HG3	2.03	0.59
18:CS:38:THR:HG22	18:CS:39:ILE:H	1.66	0.59
23:DB:1013:C:H2'	23:DB:1014:A:H8	1.66	0.59
23:DB:1447:C:H2'	23:DB:1448:G:C8	2.37	0.59
23:DB:1550:C:H2'	23:DB:1551:A:H8	1.67	0.59
23:DB:1742:U:H2'	23:DB:1743:G:C8	2.38	0.59
23:DB:2220:U:H2'	23:DB:2221:G:H8	1.68	0.59
23:DB:2324:U:H3'	23:DB:2325:G:C5'	2.32	0.59
23:DB:2852:G:H2'	23:DB:2853:C:H6	1.67	0.59
23:DB:2899:A:H2'	23:DB:2900:A:C8	2.37	0.59
26:DD:31:ALA:CB	26:DD:95:SER:HB2	2.32	0.59
26:DD:52:THR:OG1	26:DD:75:ALA:HB1	2.02	0.59
27:DE:116:ASP:CB	27:DE:185:LYS:HA	2.30	0.59
23:DB:659:G:H4'	27:DE:95:LYS:HD3	1.85	0.59
28:DF:135:ILE:O	28:DF:137:PHE:N	2.36	0.59
31:DJ:15:TRP:CD2	31:DJ:138:GLN:HB2	2.38	0.59
34:DM:131:VAL:HG22	34:DM:133:LYS:H	1.68	0.59
36:DO:74:VAL:O	36:DO:78:VAL:HG23	2.02	0.59
37:DP:90:ALA:HB3	37:DP:112:ARG:H	1.67	0.59
24:DV:9:ARG:HH21	24:DV:20:LEU:HD11	1.68	0.59
1:AA:1130:A:H2'	1:AA:1131:G:H8	1.68	0.59
1:AA:766:A:H2	1:AA:1525:G:N3	1.99	0.59
1:AA:796:C:H5'	10:AK:128:VAL:HG13	1.84	0.59
5:AF:18:VAL:O	5:AF:22:ILE:HG13	2.02	0.59
16:AQ:4:ILE:HD12	16:AQ:4:ILE:H	1.67	0.59
48:B1:42:VAL:HB	48:B1:43:ARG:HH21	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1312:U:OP2	41:BT:66:LYS:HG2	2.02	0.59
23:BB:1470:A:H3'	23:BB:1471:G:H8	1.67	0.59
23:BB:209:C:H2'	23:BB:210:C:H6	1.67	0.59
23:BB:222:A:N1	23:BB:233:A:H5''	2.17	0.59
25:BC:163:ILE:HG22	25:BC:164:VAL:N	2.16	0.59
27:BE:132:LYS:HD3	27:BE:133:LEU:CD2	2.32	0.59
27:BE:193:VAL:HG12	27:BE:194:LYS:HD2	1.84	0.59
29:BG:102:ILE:HG23	29:BG:104:LEU:HD11	1.83	0.59
29:BG:18:ILE:HD13	29:BG:18:ILE:H	1.66	0.59
52:BI:14:ALA:HA	52:BI:45:THR:HG21	1.83	0.59
34:BM:73:ILE:HG12	34:BM:90:GLU:OE2	2.03	0.59
37:BP:92:ARG:CZ	37:BP:110:LYS:HB3	2.33	0.59
38:BQ:102:LYS:HD2	38:BQ:102:LYS:N	2.18	0.59
1:CA:121:U:H3'	1:CA:121:U:OP1	2.03	0.59
3:CD:2:ARG:HH11	3:CD:114:ARG:HD3	1.65	0.59
5:CF:92:THR:HG22	5:CF:93:LYS:N	2.18	0.59
23:DB:1262:A:C2	47:D0:6:LYS:HD2	2.36	0.59
23:DB:2758:A:H2'	23:DB:2759:G:H5'	1.85	0.59
23:DB:634:C:H2'	23:DB:635:C:C6	2.38	0.59
23:DB:1825:U:H5'	25:DC:244:VAL:HG21	1.83	0.59
23:DB:587:C:H5'	27:DE:85:PHE:CE1	2.37	0.59
28:DF:26:GLN:HG2	28:DF:26:GLN:O	2.03	0.59
30:DH:103:VAL:HG13	30:DH:108:VAL:HB	1.84	0.59
30:DH:129:GLU:HA	30:DH:143:ILE:HA	1.85	0.59
31:DJ:36:LEU:HD13	31:DJ:54:ILE:HD12	1.83	0.59
36:DO:15:ARG:HH22	43:DW:76:ARG:HG2	1.68	0.59
1:AA:390:U:H2'	1:AA:391:G:H8	1.62	0.59
20:AB:206:ILE:HD13	20:AB:206:ILE:O	2.02	0.59
20:AB:46:VAL:HA	20:AB:49:PHE:HD2	1.68	0.59
2:AC:21:TRP:CB	2:AC:58:ARG:HB2	2.33	0.59
8:AI:35:GLU:HA	8:AI:39:GLY:CA	2.33	0.59
1:AA:716:A:N3	10:AK:119:GLY:HA2	2.17	0.59
10:AK:81:LEU:HD22	10:AK:104:PHE:HB3	1.84	0.59
1:AA:522:C:H41	11:AL:49:ARG:HH22	1.49	0.59
19:AT:54:GLN:N	19:AT:55:PRO:HD2	2.17	0.59
47:B0:2:VAL:HG22	47:B0:3:GLN:H	1.67	0.59
23:BB:1241:A:H2'	23:BB:1242:U:C5'	2.32	0.59
23:BB:2617:U:H2'	23:BB:2618:G:H5'	1.85	0.59
23:BB:419:U:H2'	23:BB:420:C:H6	1.67	0.59
23:BB:441:U:H2'	23:BB:442:G:H8	1.67	0.59
25:BC:30:ALA:N	25:BC:31:PRO:CD	2.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:70:ARG:HA	28:BF:80:GLN:HG3	1.83	0.59
23:BB:1007:C:H5''	31:BJ:37:ARG:NH2	2.18	0.59
33:BL:131:ALA:HB1	33:BL:135:ILE:HD12	1.85	0.59
23:BB:2467:C:O4'	34:BM:118:LYS:HG2	2.02	0.59
34:BM:86:LYS:HD2	34:BM:86:LYS:H	1.68	0.59
23:BB:483:A:H4'	42:BU:45:GLN:HA	1.83	0.59
24:BV:40:ILE:N	24:BV:40:ILE:HD13	2.18	0.59
1:CA:1031:C:H4'	1:CA:1032:G:H5''	1.84	0.59
1:CA:1326:U:O2'	1:CA:1327:C:H5'	2.02	0.59
2:CC:65:VAL:HG21	2:CC:90:VAL:HG11	1.83	0.59
21:CU:24:LYS:HD2	21:CU:25:ALA:N	2.11	0.59
23:DB:1111:A:H4'	29:DG:1:SER:O	2.03	0.59
23:DB:1210:G:H4'	23:DB:1211:C:OP2	2.02	0.59
23:DB:1412:U:H2'	23:DB:1413:A:C8	2.38	0.59
23:DB:213:A:O2'	23:DB:214:G:H5'	2.03	0.59
23:DB:2859:G:H2'	23:DB:2860:A:C8	2.36	0.59
23:DB:299:A:N6	23:DB:322:A:O2'	2.33	0.59
23:DB:324:A:OP2	27:DE:164:LEU:HB2	2.02	0.59
23:DB:79:C:O2'	23:DB:346:A:H1'	2.03	0.59
25:DC:163:ILE:HG12	25:DC:173:LEU:CD2	2.33	0.59
25:DC:261:ARG:HG2	25:DC:261:ARG:HH11	1.66	0.59
26:DD:22:ILE:HG13	26:DD:22:ILE:O	2.01	0.59
26:DD:77:ARG:HB2	26:DD:77:ARG:HH11	1.67	0.59
28:DF:133:GLU:H	28:DF:150:GLY:HA2	1.66	0.59
30:DH:89:LYS:HA	30:DH:123:ARG:O	2.03	0.59
52:DI:17:ALA:C	52:DI:19:PRO:HD3	2.23	0.59
23:DB:1099:G:O4'	52:DI:3:LYS:O	2.20	0.59
33:DL:2:ARG:HG3	33:DL:2:ARG:NH1	2.18	0.59
33:DL:64:PHE:HD2	50:D3:11:LYS:HB3	1.68	0.59
33:DL:85:VAL:HG13	33:DL:86:GLU:N	2.11	0.59
34:DM:40:ARG:HA	34:DM:92:TRP:NE1	2.17	0.59
35:DN:41:ALA:HB1	35:DN:113:ILE:CD1	2.32	0.59
24:DV:9:ARG:HG2	24:DV:41:GLU:CG	2.32	0.59
44:DX:25:GLN:O	44:DX:29:ARG:HG3	2.02	0.59
1:AA:121:U:H3'	1:AA:121:U:OP1	2.03	0.59
1:AA:801:U:H2'	1:AA:802:A:H8	1.67	0.59
4:AE:110:MET:O	4:AE:113:VAL:HG22	2.03	0.59
6:AG:117:LEU:HD22	6:AG:120:ALA:HB3	1.84	0.59
12:AM:64:VAL:HG12	12:AM:65:GLU:N	2.18	0.59
23:BB:1429:G:H2'	23:BB:1430:G:C8	2.37	0.59
23:BB:156:A:H2'	23:BB:157:C:H6	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2097:A:H2'	23:BB:2098:U:C6	2.38	0.59
23:BB:2328:A:H2'	23:BB:2329:U:H6	1.67	0.59
23:BB:296:U:H2'	23:BB:297:G:H8	1.67	0.59
23:BB:355:U:H2'	23:BB:356:G:H8	1.68	0.59
23:BB:544:C:H2'	23:BB:545:U:C2	2.38	0.59
30:BH:62:LEU:O	30:BH:66:ASN:HB3	2.03	0.59
52:BI:32:VAL:HG22	52:BI:60:VAL:HG21	1.85	0.59
35:BN:24:MET:O	35:BN:27:SER:HB3	2.03	0.59
39:BR:47:VAL:HG13	39:BR:49:ILE:N	2.13	0.59
46:BZ:49:ARG:C	46:BZ:51:VAL:H	2.07	0.59
1:CA:1238:A:C2	1:CA:1241:G:H1'	2.38	0.59
1:CA:501:C:H1'	1:CA:549:C:H1'	1.85	0.59
1:CA:812:G:O2'	1:CA:813:U:H6	1.85	0.59
1:CA:993:G:H2'	1:CA:995:C:H41	1.66	0.59
2:CC:61:LYS:HZ3	2:CC:96:VAL:HG12	1.68	0.59
1:CA:939:G:H5''	6:CG:101:ARG:NH2	2.17	0.59
13:CN:86:ALA:HB1	13:CN:91:GLU:HB2	1.84	0.59
23:DB:2181:U:H6	23:DB:2181:U:O5'	1.86	0.59
23:DB:2369:A:O2'	23:DB:2370:G:H5'	2.03	0.59
23:DB:635:C:H3'	33:DL:126:ARG:NH2	2.18	0.59
23:DB:850:U:O2'	45:DY:22:THR:HG22	2.03	0.59
25:DC:140:VAL:HG11	25:DC:163:ILE:HD11	1.84	0.59
25:DC:171:VAL:HB	25:DC:182:LYS:CB	2.26	0.59
52:DI:2:LYS:C	52:DI:3:LYS:HD2	2.23	0.59
32:DK:99:ILE:CD1	32:DK:115:ILE:HG13	2.32	0.59
33:DL:109:LYS:CG	33:DL:126:ARG:HD3	2.33	0.59
39:DR:6:GLN:OE1	39:DR:38:VAL:HG22	2.02	0.59
40:DS:27:LYS:HG3	40:DS:70:LYS:HD3	1.85	0.59
23:DB:95:A:O2'	44:DX:43:LEU:HD21	2.03	0.59
1:AA:202:G:H2'	1:AA:203:G:C8	2.37	0.59
1:AA:300:A:H2'	1:AA:301:G:O4'	2.03	0.59
1:AA:1348:U:OP1	8:AI:111:GLU:HB2	2.02	0.59
16:AQ:26:ARG:HH21	16:AQ:39:ARG:NH1	2.00	0.59
18:AS:11:ASP:CG	18:AS:34:SER:HB2	2.23	0.59
23:BB:1579:A:H2'	23:BB:1580:A:C8	2.38	0.59
23:BB:170:U:H2'	23:BB:171:U:C6	2.37	0.59
23:BB:2188:U:H2'	23:BB:2189:U:C6	2.37	0.59
23:BB:2491:U:H5''	23:BB:2570:G:C5'	2.33	0.59
23:BB:2789:C:H3'	23:BB:2893:A:H62	1.66	0.59
28:BF:15:LEU:HD13	28:BF:28:PRO:HG2	1.83	0.59
29:BG:17:LYS:HB3	29:BG:24:THR:HG22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BI:18:ASN:N	52:BI:19:PRO:HD2	2.17	0.59
52:BI:77:VAL:HA	52:BI:80:LYS:CE	2.33	0.59
23:BB:1076:C:H4'	52:BI:94:LYS:CE	2.33	0.59
35:BN:38:LEU:HB2	35:BN:109:PRO:CB	2.33	0.59
37:BP:59:THR:HA	37:BP:76:HIS:HA	1.84	0.59
39:BR:90:ARG:HH11	39:BR:91:GLN:N	2.01	0.59
40:BS:66:ILE:N	40:BS:66:ILE:HD12	2.18	0.59
42:BU:13:LEU:HD12	42:BU:69:VAL:HA	1.84	0.59
42:BU:73:ASN:HB2	42:BU:96:LYS:NZ	2.18	0.59
23:BB:2261:C:OP2	43:BW:13:ARG:HB2	2.03	0.59
44:BX:35:GLY:CA	44:BX:40:SER:HA	2.32	0.59
46:BZ:15:SER:O	46:BZ:21:VAL:HB	2.02	0.59
1:CA:1019:A:H2'	1:CA:1020:G:C8	2.38	0.59
1:CA:131:A:H2'	1:CA:132:C:H6	1.68	0.59
20:CB:132:GLU:HA	20:CB:135:MET:HE2	1.85	0.59
4:CE:95:MET:HG3	4:CE:124:ALA:HB2	1.83	0.59
7:CH:54:THR:HG23	7:CH:55:LYS:HG2	1.83	0.59
8:CI:38:PHE:O	8:CI:44:ARG:HD2	2.03	0.59
1:CA:529:G:O6	11:CL:45:ASN:HA	2.03	0.59
14:CO:28:VAL:HG11	14:CO:80:LEU:HD21	1.85	0.59
49:D2:13:ASN:HA	49:D2:16:HIS:O	2.02	0.59
23:DB:1442:U:H2'	23:DB:1443:U:H6	1.68	0.59
23:DB:1470:A:H3'	23:DB:1471:G:H8	1.66	0.59
23:DB:794:A:H2'	23:DB:795:C:H6	1.67	0.59
23:DB:873:C:H2'	23:DB:874:G:H8	1.68	0.59
28:DF:65:LEU:CD2	28:DF:87:LYS:HD2	2.33	0.59
31:DJ:21:THR:HG22	31:DJ:58:ASN:OD1	2.03	0.59
42:DU:31:GLY:O	42:DU:32:LYS:HD3	2.03	0.59
1:AA:1448:C:H2'	1:AA:1449:C:C6	2.38	0.58
1:AA:403:C:O2'	1:AA:404:G:H5'	2.02	0.58
3:AD:144:ILE:CD1	3:AD:154:VAL:HG21	2.33	0.58
12:AM:2:ARG:HB3	12:AM:6:ILE:O	2.03	0.58
19:AT:53:MET:O	19:AT:57:VAL:HG23	2.02	0.58
50:B3:23:HIS:HD2	50:B3:23:HIS:H	1.50	0.58
23:BB:131:A:H2'	23:BB:132:G:C8	2.38	0.58
23:BB:1387:A:H5'	23:BB:1469:A:H1'	1.84	0.58
23:BB:1454:C:H5'	35:BN:63:ARG:NH2	2.07	0.58
23:BB:1484:U:H2'	23:BB:1485:U:C6	2.37	0.58
23:BB:1484:U:H2'	23:BB:1485:U:H6	1.68	0.58
23:BB:1652:A:H62	35:BN:11:ASN:ND2	2.01	0.58
23:BB:2789:C:H2'	23:BB:2893:A:N7	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:126:ASN:N	28:BF:126:ASN:ND2	2.50	0.58
31:BJ:20:ALA:HB1	31:BJ:59:ALA:CA	2.33	0.58
32:BK:104:THR:O	32:BK:107:LEU:HB3	2.03	0.58
32:BK:60:ALA:HA	32:BK:87:LEU:HG	1.84	0.58
38:BQ:97:ILE:HA	38:BQ:100:PHE:CZ	2.38	0.58
40:BS:7:HIS:CD2	40:BS:10:ALA:HB2	2.38	0.58
41:BT:36:LYS:O	41:BT:36:LYS:HD3	2.01	0.58
1:CA:34:C:H2'	1:CA:35:G:H8	1.68	0.58
1:CA:476:U:H2'	1:CA:477:C:C6	2.38	0.58
4:CE:24:VAL:HG23	4:CE:26:GLY:H	1.67	0.58
4:CE:61:LYS:O	4:CE:64:GLU:HB3	2.03	0.58
23:DB:336:C:H5''	42:DU:3:LYS:HZ1	1.68	0.58
23:DB:672:C:O2'	23:DB:673:C:H5'	2.03	0.58
25:DC:237:ARG:C	25:DC:237:ARG:HE	2.07	0.58
26:DD:110:THR:HB	26:DD:202:ILE:HB	1.84	0.58
27:DE:152:GLU:O	27:DE:153:LEU:HB2	2.02	0.58
27:DE:189:THR:C	27:DE:191:ASP:H	2.06	0.58
30:DH:98:ASP:HA	30:DH:101:ASP:OD1	2.03	0.58
32:DK:24:VAL:HG13	32:DK:33:ALA:HB2	1.85	0.58
33:DL:118:THR:OG1	33:DL:138:ALA:HB2	2.02	0.58
33:DL:123:ARG:HB3	33:DL:141:LYS:HB3	1.83	0.58
36:DO:93:ASP:HB2	36:DO:117:PHE:CE2	2.38	0.58
37:DP:65:ASN:H	37:DP:71:ARG:HA	1.68	0.58
37:DP:25:VAL:CG1	37:DP:87:ARG:HA	2.29	0.58
38:DQ:20:ALA:HA	38:DQ:23:TYR:CE2	2.37	0.58
38:DQ:87:VAL:HB	39:DR:54:VAL:HG11	1.85	0.58
23:DB:136:G:N1	41:DT:3:ARG:NH1	2.51	0.58
42:DU:98:ASN:O	42:DU:99:SER:HB2	2.01	0.58
43:DW:65:LYS:HB2	43:DW:65:LYS:NZ	2.18	0.58
1:AA:1011:C:H2'	1:AA:1012:A:C8	2.38	0.58
1:AA:1237:C:H3'	1:AA:1238:A:H5'	1.85	0.58
1:AA:241:G:O2'	1:AA:242:G:H5'	2.03	0.58
1:AA:687:A:C2	1:AA:704:A:C5	2.91	0.58
20:AB:60:ALA:HB3	20:AB:223:GLY:HA3	1.83	0.58
1:AA:408:A:H5'	3:AD:21:LYS:HE2	1.84	0.58
5:AF:42:TRP:HB2	5:AF:59:TYR:HB2	1.84	0.58
12:AM:72:ILE:O	12:AM:76:ILE:HG13	2.03	0.58
15:AP:46:LYS:HE2	15:AP:46:LYS:H	1.65	0.58
10:AK:113:THR:CG2	21:AU:28:LEU:HD21	2.33	0.58
23:BB:156:A:H2'	23:BB:157:C:C6	2.38	0.58
23:BB:1945:G:H2'	23:BB:1946:U:C6	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2369:A:O2'	23:BB:2370:G:H5'	2.03	0.58
23:BB:2386:A:H2'	23:BB:2387:U:C6	2.38	0.58
23:BB:2480:C:O2'	23:BB:2481:G:H5'	2.03	0.58
23:BB:686:U:O4	49:B2:12:ARG:HG3	2.02	0.58
30:BH:6:LEU:HA	30:BH:15:LEU:C	2.23	0.58
30:BH:94:ILE:HG23	30:BH:99:ILE:HD11	1.84	0.58
52:BI:105:LEU:HD11	52:BI:139:VAL:CG1	2.33	0.58
31:BJ:18:VAL:HG22	31:BJ:55:ILE:O	2.02	0.58
31:BJ:77:HIS:CD2	31:BJ:84:ILE:HB	2.38	0.58
34:BM:38:ARG:HH11	34:BM:38:ARG:CB	2.10	0.58
37:BP:28:LYS:HG2	37:BP:44:GLY:O	2.03	0.58
37:BP:31:VAL:HG12	37:BP:33:GLU:H	1.68	0.58
37:BP:59:THR:HG23	37:BP:76:HIS:NE2	2.18	0.58
42:BU:70:ALA:HB3	42:BU:77:GLY:O	2.03	0.58
24:BV:84:PRO:HG3	34:BM:127:LYS:HE2	1.85	0.58
1:CA:1119:C:H2'	1:CA:1120:C:C6	2.37	0.58
1:CA:413:G:H4'	1:CA:414:A:H5''	1.85	0.58
3:CD:30:LYS:HB2	3:CD:30:LYS:NZ	2.18	0.58
3:CD:71:PHE:HE2	3:CD:89:LEU:HD11	1.68	0.58
12:CM:89:ARG:HB3	12:CM:96:VAL:HG22	1.84	0.58
50:D3:14:LYS:HA	50:D3:63:TYR:CE2	2.38	0.58
23:DB:1857:G:H1'	23:DB:1885:A:N6	2.16	0.58
23:DB:2022:U:O2'	23:DB:2617:U:H5'	2.03	0.58
23:DB:2836:U:H2'	23:DB:2837:A:C8	2.39	0.58
23:DB:2887:A:N9	47:D0:27:LEU:HD21	2.18	0.58
23:DB:28:A:O2'	23:DB:29:U:H5'	2.03	0.58
25:DC:87:SER:N	25:DC:155:ARG:HH12	2.01	0.58
28:DF:107:VAL:HG11	28:DF:175:PRO:HG3	1.85	0.58
29:DG:41:GLU:HG2	29:DG:54:ARG:HH21	1.68	0.58
34:DM:41:LEU:HG	34:DM:46:ILE:HD11	1.85	0.58
37:DP:7:LEU:O	37:DP:11:GLN:HG2	2.03	0.58
38:DQ:39:ILE:HG13	38:DQ:40:LYS:N	2.17	0.58
1:AA:1246:A:H2'	1:AA:1247:U:C6	2.38	0.58
1:AA:203:G:N2	1:AA:205:A:H61	2.01	0.58
1:AA:950:U:H2'	1:AA:951:G:H8	1.68	0.58
20:AB:67:LEU:HD13	20:AB:67:LEU:N	2.18	0.58
13:AN:71:GLY:O	13:AN:79:SER:HA	2.04	0.58
15:AP:26:ASN:CG	15:AP:31:ARG:HB3	2.24	0.58
18:AS:4:LEU:HD13	18:AS:8:PRO:HA	1.84	0.58
51:B4:16:ILE:N	51:B4:16:ILE:HD12	2.17	0.58
23:BB:1447:C:H2'	23:BB:1448:G:C8	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1552:A:C2'	23:BB:1553:A:H5'	2.33	0.58
23:BB:2836:U:H2'	23:BB:2837:A:C8	2.39	0.58
23:BB:877:A:N1	23:BB:899:A:H2'	2.17	0.58
25:BC:32:LEU:HA	25:BC:81:GLU:OE1	2.04	0.58
27:BE:3:LEU:CB	27:BE:12:LEU:HD22	2.33	0.58
23:BB:37:C:H2'	27:BE:46:GLN:OE1	2.03	0.58
28:BF:102:LEU:C	28:BF:107:VAL:HG23	2.23	0.58
28:BF:132:ARG:NE	28:BF:147:ARG:HE	2.00	0.58
30:BH:35:LYS:O	30:BH:37:VAL:HG13	2.03	0.58
30:BH:73:ASN:ND2	30:BH:73:ASN:N	2.52	0.58
31:BJ:75:TYR:HA	31:BJ:86:GLN:O	2.04	0.58
32:BK:11:ALA:HB3	32:BK:85:VAL:HG23	1.85	0.58
42:BU:46:LYS:HE2	42:BU:54:PRO:O	2.03	0.58
1:CA:1163:A:H2'	1:CA:1164:G:H8	1.67	0.58
1:CA:300:A:H2'	1:CA:301:G:O4'	2.03	0.58
4:CE:113:VAL:HG23	4:CE:114:LEU:N	2.18	0.58
4:CE:11:GLN:OE1	4:CE:116:VAL:HA	2.03	0.58
8:CI:41:GLU:H	8:CI:44:ARG:HD3	1.68	0.58
9:CJ:38:GLY:O	9:CJ:74:VAL:HA	2.03	0.58
9:CJ:85:ASP:HA	9:CJ:89:ARG:HB2	1.85	0.58
10:CK:73:VAL:O	10:CK:76:TYR:HB2	2.04	0.58
14:CO:20:ASP:OD2	14:CO:23:SER:HB2	2.04	0.58
23:DB:1229:C:H2'	23:DB:1230:A:H8	1.68	0.58
23:DB:1515:A:H4'	23:DB:1556:C:O2'	2.04	0.58
23:DB:230:G:H2'	23:DB:231:A:H8	1.67	0.58
23:DB:2783:U:H2'	23:DB:2784:U:H6	1.68	0.58
23:DB:307:G:N2	23:DB:309:A:H3'	2.17	0.58
23:DB:419:U:H2'	23:DB:420:C:C6	2.38	0.58
23:DB:764:A:H5'	25:DC:208:GLY:HA3	1.84	0.58
23:DB:946:C:H2'	23:DB:947:A:H8	1.68	0.58
25:DC:109:LEU:HD21	25:DC:115:ILE:HD11	1.84	0.58
25:DC:86:ARG:C	25:DC:155:ARG:HH12	2.06	0.58
27:DE:14:VAL:HG12	27:DE:16:GLU:H	1.67	0.58
30:DH:5:LEU:HD23	30:DH:17:ASP:O	2.02	0.58
30:DH:3:VAL:O	30:DH:18:GLN:HA	2.03	0.58
52:DI:85:ILE:CD1	52:DI:137:LEU:HD21	2.32	0.58
33:DL:7:SER:HB2	33:DL:8:PRO:CD	2.17	0.58
22:DA:38:C:H4'	36:DO:100:HIS:NE2	2.18	0.58
37:DP:38:ARG:CG	37:DP:39:LEU:H	2.08	0.58
38:DQ:39:ILE:O	38:DQ:43:GLN:HB3	2.02	0.58
38:DQ:89:ILE:H	38:DQ:89:ILE:HD12	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DR:71:LYS:HB2	39:DR:73:LYS:NZ	2.18	0.58
42:DU:11:ILE:HB	42:DU:69:VAL:HG23	1.84	0.58
1:AA:1324:A:H2'	1:AA:1325:C:H6	1.69	0.58
1:AA:621:A:H2'	1:AA:622:A:C8	2.38	0.58
1:AA:88:U:O2'	1:AA:89:U:C6	2.56	0.58
2:AC:151:GLU:O	2:AC:197:VAL:HA	2.02	0.58
10:AK:26:PHE:HE1	10:AK:88:PRO:HG2	1.68	0.58
11:AL:22:ALA:HB2	11:AL:56:LEU:HD21	1.85	0.58
16:AQ:66:LEU:HD11	16:AQ:73:THR:HG22	1.85	0.58
51:B4:16:ILE:H	51:B4:16:ILE:HD12	1.68	0.58
23:BB:1082:U:N3	23:BB:1086:A:C2	2.71	0.58
23:BB:111:A:H2'	23:BB:112:U:O4'	2.02	0.58
23:BB:1144:A:O2'	23:BB:1145:C:H5'	2.03	0.58
23:BB:1229:C:H2'	23:BB:1230:A:C8	2.38	0.58
23:BB:2898:U:H2'	23:BB:2899:A:H8	1.69	0.58
23:BB:27:G:N2	23:BB:512:G:H2'	2.17	0.58
23:BB:794:A:H2'	23:BB:795:C:H6	1.67	0.58
25:BC:119:VAL:HG13	25:BC:133:ASN:ND2	2.18	0.58
25:BC:181:ARG:NH1	25:BC:260:LYS:HD2	2.18	0.58
25:BC:140:VAL:C	25:BC:193:GLU:HB2	2.23	0.58
26:BD:83:ARG:HH12	26:BD:85:ALA:HB2	1.67	0.58
40:BS:50:VAL:HG11	40:BS:103:ILE:HG12	1.85	0.58
43:BW:40:ARG:HH12	43:BW:68:PHE:HA	1.67	0.58
1:CA:241:G:O2'	1:CA:242:G:H5'	2.02	0.58
1:CA:620:C:C2	3:CD:131:ILE:HD13	2.38	0.58
4:CE:104:ILE:O	4:CE:104:ILE:HG23	2.03	0.58
6:CG:37:THR:O	6:CG:41:ILE:HG13	2.03	0.58
8:CI:50:PRO:HD3	8:CI:79:ARG:HD3	1.85	0.58
10:CK:28:ASN:ND2	10:CK:29:THR:H	2.00	0.58
23:DB:125:A:H4'	49:D2:13:ASN:ND2	2.18	0.58
23:DB:1579:A:H2'	23:DB:1580:A:C8	2.38	0.58
23:DB:2756:U:C1'	23:DB:2757:A:H5''	2.34	0.58
23:DB:69:C:H2'	23:DB:70:G:H8	1.68	0.58
23:DB:90:U:H2'	23:DB:91:A:C2	2.39	0.58
23:DB:942:G:O2'	23:DB:943:A:H5'	2.03	0.58
25:DC:28:PRO:HG2	25:DC:79:ARG:NH2	2.18	0.58
26:DD:24:VAL:CG1	26:DD:193:VAL:HG21	2.34	0.58
29:DG:156:TYR:HA	29:DG:171:LYS:HG3	1.84	0.58
29:DG:70:LEU:O	29:DG:74:MET:HG3	2.03	0.58
30:DH:115:VAL:HB	30:DH:132:PHE:HD1	1.68	0.58
30:DH:135:HIS:HB3	30:DH:138:VAL:CG2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:16:ARG:NH2	34:DM:72:PRO:HG2	2.17	0.58
38:DQ:15:LYS:O	38:DQ:19:GLN:HG2	2.03	0.58
39:DR:6:GLN:HB3	39:DR:41:ILE:HD13	1.84	0.58
42:DU:32:LYS:HA	42:DU:65:GLN:HE21	1.68	0.58
24:DV:43:ASP:OD2	24:DV:46:LYS:HB2	2.03	0.58
1:AA:186:C:H4'	19:AT:75:LYS:HB2	1.85	0.58
1:AA:476:U:H2'	1:AA:477:C:C6	2.38	0.58
20:AB:172:ILE:HG22	20:AB:176:ASN:HD21	1.68	0.58
2:AC:5:HIS:CD2	2:AC:7:ASN:HB2	2.38	0.58
33:BL:64:PHE:CD1	50:B3:24:LYS:HE3	2.39	0.58
23:BB:1259:G:H2'	23:BB:1260:A:H8	1.68	0.58
23:BB:1537:G:H3'	23:BB:1537:G:N3	2.18	0.58
23:BB:172:A:H2'	23:BB:173:A:H8	1.68	0.58
23:BB:1777:U:O2'	23:BB:1778:U:H5'	2.03	0.58
23:BB:2267:A:N6	23:BB:2272:U:C4	2.67	0.58
23:BB:2491:U:H5''	23:BB:2570:G:H5''	1.85	0.58
23:BB:2873:A:N3	35:BN:6:SER:HA	2.18	0.58
23:BB:917:A:C2	23:BB:918:A:H1'	2.38	0.58
23:BB:2811:G:OP1	26:BD:61:THR:HB	2.04	0.58
28:BF:107:VAL:N	28:BF:108:PRO:CD	2.65	0.58
33:BL:84:LYS:O	33:BL:85:VAL:HG12	2.03	0.58
41:BT:25:GLU:HG3	41:BT:26:LYS:H	1.67	0.58
42:BU:71:ILE:HD13	42:BU:72:PHE:O	2.03	0.58
24:BV:42:LEU:HB2	24:BV:47:VAL:HG21	1.85	0.58
1:CA:1021:A:H2'	1:CA:1022:A:C8	2.38	0.58
11:CL:77:SER:O	11:CL:79:ILE:HG23	2.03	0.58
13:CN:62:ARG:N	13:CN:72:PHE:HZ	2.02	0.58
14:CO:62:ARG:CZ	14:CO:86:LEU:HD21	2.33	0.58
16:CQ:42:LYS:C	16:CQ:43:LEU:HD12	2.23	0.58
51:D4:10:LEU:HB2	51:D4:25:VAL:CG2	2.34	0.58
23:DB:2010:G:H5''	40:DS:42:LYS:HB2	1.84	0.58
23:DB:208:C:H2'	23:DB:209:C:H6	1.68	0.58
23:DB:2421:G:OP2	50:D3:31:ILE:HG21	2.04	0.58
23:DB:2491:U:H5''	23:DB:2570:G:H5''	1.85	0.58
23:DB:2893:A:H5''	23:DB:2894:G:H5'	1.85	0.58
25:DC:110:LYS:HE3	25:DC:110:LYS:HA	1.84	0.58
27:DE:147:LEU:H	27:DE:147:LEU:HD12	1.68	0.58
27:DE:151:GLY:O	27:DE:171:ASP:HA	2.04	0.58
27:DE:187:VAL:HG23	27:DE:188:MET:H	1.68	0.58
32:DK:43:ILE:N	32:DK:43:ILE:HD12	2.18	0.58
32:DK:6:THR:HG22	32:DK:7:MET:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:87:LEU:HD22	41:DT:93:LEU:HD11	1.85	0.58
44:DX:56:LEU:HA	44:DX:59:GLU:HB2	1.86	0.58
1:AA:1366:C:H2'	1:AA:1367:C:H6	1.68	0.58
1:AA:922:G:N3	1:AA:1398:A:H2	2.02	0.58
1:AA:1508:A:H2'	1:AA:1509:C:C6	2.38	0.58
1:AA:923:A:H2'	1:AA:924:C:C6	2.39	0.58
14:AO:68:TYR:CZ	14:AO:72:LYS:HG3	2.39	0.58
50:B3:12:ARG:HG2	50:B3:23:HIS:HB3	1.85	0.58
23:BB:1046:A:C3'	23:BB:1047:G:H5''	2.32	0.58
23:BB:1709:U:H2'	23:BB:1710:G:H8	1.68	0.58
23:BB:1742:U:H2'	23:BB:1743:G:C8	2.38	0.58
23:BB:1799:G:C5	25:BC:175:LEU:HD13	2.39	0.58
23:BB:2083:G:H2'	23:BB:2084:C:C6	2.38	0.58
23:BB:26:G:H2'	23:BB:27:G:C1'	2.34	0.58
23:BB:38:A:C8	27:BE:46:GLN:OE1	2.57	0.58
23:BB:950:G:H2'	23:BB:951:C:C6	2.39	0.58
26:BD:14:ILE:HA	26:BD:19:GLY:HA2	1.85	0.58
30:BH:57:LYS:HG3	30:BH:61:VAL:CG2	2.34	0.58
33:BL:78:ARG:CZ	33:BL:110:VAL:HG11	2.33	0.58
33:BL:54:GLN:HB3	33:BL:56:PRO:HD2	1.84	0.58
34:BM:101:VAL:C	34:BM:102:LEU:HD12	2.24	0.58
34:BM:71:LYS:HE2	34:BM:91:TYR:C	2.23	0.58
39:BR:20:VAL:HG13	39:BR:97:LYS:HD2	1.86	0.58
1:CA:1320:C:H41	18:CS:36:ARG:HD2	1.68	0.58
1:CA:202:G:H2'	1:CA:203:G:H8	1.69	0.58
23:DB:1444:G:H2'	23:DB:1445:G:H8	1.67	0.58
23:DB:1484:U:H2'	23:DB:1485:U:H6	1.68	0.58
23:DB:1557:C:H3'	23:DB:1558:C:H5''	1.84	0.58
23:DB:1777:U:O2'	23:DB:1778:U:H5'	2.03	0.58
23:DB:2491:U:H5''	23:DB:2570:G:C5'	2.33	0.58
23:DB:98:G:C3'	23:DB:99:U:H5''	2.33	0.58
31:DJ:94:ALA:HB1	31:DJ:95:ARG:HH21	1.69	0.58
33:DL:25:SER:C	33:DL:27:LEU:H	2.07	0.58
37:DP:71:ARG:HH11	37:DP:71:ARG:HB3	1.67	0.58
1:AA:1100:C:OP2	20:AB:94:ARG:HD3	2.04	0.58
1:AA:1130:A:O2'	1:AA:1131:G:H5'	2.04	0.58
1:AA:398:U:H2'	1:AA:399:G:C8	2.39	0.58
2:AC:71:ARG:NH2	2:AC:74:ILE:HB	2.18	0.58
3:AD:17:ASP:OD2	3:AD:27:ILE:HG12	2.04	0.58
3:AD:60:VAL:HB	3:AD:194:ILE:HG13	1.84	0.58
4:AE:158:LYS:N	4:AE:158:LYS:HE2	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AJ:71:LEU:H	9:AJ:71:LEU:HD12	1.68	0.58
48:B1:35:LEU:C	48:B1:36:LYS:HD2	2.24	0.58
23:BB:2351:G:O6	50:B3:38:LYS:HD2	2.03	0.58
23:BB:146:A:H2'	23:BB:147:C:H6	1.66	0.58
23:BB:1847:A:H4'	23:BB:1848:A:C8	2.39	0.58
23:BB:1932:A:H2'	23:BB:1933:G:O4'	2.04	0.58
23:BB:19:A:H2'	23:BB:20:C:H6	1.65	0.58
23:BB:21:A:O2'	23:BB:22:C:H5'	2.04	0.58
23:BB:3:U:O2'	23:BB:4:U:H6	1.87	0.58
25:BC:75:ALA:HB2	25:BC:95:TYR:CE1	2.38	0.58
26:BD:138:LEU:HD23	26:BD:139:SER:H	1.68	0.58
33:BL:133:ALA:HA	33:BL:136:GLU:OE2	2.03	0.58
34:BM:67:VAL:HB	34:BM:100:LYS:CG	2.32	0.58
35:BN:29:VAL:HG22	35:BN:83:LEU:HD22	1.85	0.58
36:BO:21:LEU:HD13	36:BO:22:GLY:N	2.18	0.58
36:BO:40:ILE:HG12	36:BO:41:ALA:O	2.04	0.58
1:CA:1142:G:H2'	1:CA:1143:G:H8	1.68	0.58
1:CA:1299:A:N7	1:CA:1301:U:H1'	2.19	0.58
1:CA:212:G:H2'	1:CA:213:G:H8	1.68	0.58
1:CA:26:A:N6	1:CA:558:G:H1'	2.19	0.58
12:CM:79:LEU:HA	12:CM:82:LEU:HG	1.85	0.58
23:DB:2466:C:OP1	51:D4:5:ALA:HB3	2.04	0.58
22:DA:28:C:OP1	36:DO:33:ARG:HD3	2.04	0.58
22:DA:35:C:H2'	22:DA:36:C:O4'	2.03	0.58
23:DB:976:G:H4'	23:DB:1156:A:N7	2.19	0.58
23:DB:138:U:H2'	23:DB:140:C:C6	2.38	0.58
23:DB:1439:A:N6	23:DB:1440:U:O2	2.35	0.58
23:DB:2020:A:H5'	47:D0:8:THR:HG21	1.84	0.58
23:DB:479:A:N3	23:DB:481:G:H5''	2.19	0.58
25:DC:74:PRO:HG3	25:DC:116:GLN:HG2	1.85	0.58
23:DB:2024:G:H5''	26:DD:154:LYS:NZ	2.18	0.58
28:DF:56:LEU:HA	28:DF:59:ILE:CG2	2.34	0.58
33:DL:90:VAL:HG13	33:DL:122:VAL:HG13	1.86	0.58
41:DT:34:VAL:HG22	41:DT:35:ALA:H	1.67	0.58
41:DT:66:LYS:HA	41:DT:76:ARG:O	2.03	0.58
1:AA:1160:G:H2'	1:AA:1161:C:H6	1.69	0.58
1:AA:1299:A:H62	1:AA:1302:C:H5	1.52	0.58
1:AA:742:G:O2'	1:AA:743:A:H5'	2.03	0.58
20:AB:65:LYS:HD2	20:AB:157:PRO:HA	1.86	0.58
20:AB:166:ASP:OD1	20:AB:190:SER:HA	2.03	0.58
20:AB:187:ASP:HB2	20:AB:203:ASP:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:113:LYS:HD2	2:AC:184:ASN:OD1	2.04	0.58
3:AD:191:SER:O	3:AD:192:ALA:HB2	2.04	0.58
4:AE:140:ILE:HA	4:AE:143:LEU:HD12	1.85	0.58
10:AK:110:THR:HG22	21:AU:4:LYS:HA	1.86	0.58
23:BB:55:G:P	49:B2:35:ARG:HH21	2.26	0.58
23:BB:1199:U:H2'	23:BB:1200:C:H6	1.67	0.58
23:BB:1791:A:C2	23:BB:1829:A:C4'	2.87	0.58
23:BB:17:G:H2'	23:BB:18:U:C6	2.39	0.58
23:BB:782:A:O2'	25:BC:223:ALA:HB1	2.04	0.58
23:BB:950:G:H2'	23:BB:951:C:H6	1.67	0.58
23:BB:962:G:N2	34:BM:81:ARG:HH21	2.01	0.58
29:BG:60:GLY:O	29:BG:63:GLN:HG2	2.04	0.58
26:BD:157:LYS:HG2	31:BJ:79:GLY:O	2.02	0.58
32:BK:40:LYS:NZ	32:BK:89:ASN:HD21	2.01	0.58
33:BL:57:LEU:HD12	33:BL:57:LEU:O	2.03	0.58
23:BB:2874:C:P	35:BN:5:LYS:HG2	2.44	0.58
37:BP:7:LEU:HD13	37:BP:11:GLN:HG2	1.84	0.58
40:BS:29:VAL:CG2	40:BS:71:VAL:HG23	2.34	0.58
45:BY:22:THR:HG21	45:BY:50:VAL:HG11	1.85	0.58
46:BZ:5:ILE:CD1	46:BZ:5:ILE:H	2.11	0.58
1:CA:1219:A:H2'	1:CA:1220:G:C8	2.39	0.58
1:CA:1323:G:H2'	1:CA:1324:A:H8	1.68	0.58
1:CA:450:G:N7	1:CA:481:G:O6	2.37	0.58
3:CD:19:PHE:O	3:CD:22:SER:HB2	2.03	0.58
1:CA:797:C:OP1	10:CK:125:LYS:HG3	2.04	0.58
11:CL:80:LEU:HD23	11:CL:97:VAL:HG21	1.86	0.58
16:CQ:10:ARG:NH2	16:CQ:55:GLY:N	2.52	0.58
17:CR:23:LYS:HD2	17:CR:23:LYS:O	2.04	0.58
47:D0:27:LEU:H	47:D0:27:LEU:CD2	2.15	0.58
50:D3:12:ARG:NE	50:D3:23:HIS:HB2	2.19	0.58
22:DA:102:G:H2'	22:DA:103:U:C6	2.39	0.58
25:DC:21:PRO:HG2	25:DC:22:GLU:H	1.68	0.58
26:DD:148:GLN:HB3	26:DD:151:THR:CG2	2.34	0.58
27:DE:115:GLN:CG	27:DE:116:ASP:H	2.17	0.58
27:DE:152:GLU:HB2	27:DE:158:PHE:CE1	2.38	0.58
30:DH:94:ILE:CG2	30:DH:122:LEU:HG	2.33	0.58
52:DI:76:ALA:HA	52:DI:135:MET:SD	2.44	0.58
34:DM:81:ARG:HG3	34:DM:82:MET:HG2	1.85	0.58
35:DN:35:LYS:HD3	35:DN:112:TYR:CD1	2.39	0.58
38:DQ:109:VAL:O	38:DQ:113:LYS:HB2	2.04	0.58
40:DS:2:GLU:HB2	40:DS:108:SER:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DS:39:THR:O	40:DS:39:THR:HG23	2.03	0.58
42:DU:8:ASP:O	42:DU:10:VAL:N	2.37	0.58
43:DW:44:PHE:CE2	43:DW:60:ALA:HB2	2.38	0.58
45:DY:38:GLU:O	45:DY:43:ILE:HG21	2.04	0.58
46:DZ:11:GLU:H	46:DZ:27:THR:CG2	2.16	0.58
1:AA:1171:A:H2'	1:AA:1172:C:C6	2.38	0.58
1:AA:1468:A:H2'	1:AA:1469:C:O4'	2.03	0.58
1:AA:1519:A:H3'	1:AA:1520:C:C5'	2.34	0.58
1:AA:26:A:N6	1:AA:558:G:H1'	2.19	0.58
1:AA:80:A:H2'	1:AA:81:A:O4'	2.04	0.58
3:AD:138:PRO:HA	3:AD:181:PHE:CD2	2.39	0.58
4:AE:111:ARG:O	4:AE:115:GLU:HG3	2.04	0.58
4:AE:13:LYS:HE3	4:AE:115:GLU:HB2	1.86	0.58
5:AF:64:VAL:HG12	5:AF:65:GLU:N	2.15	0.58
8:AI:51:LEU:HB3	8:AI:56:MET:HG2	1.85	0.58
9:AJ:40:ILE:HG22	9:AJ:41:PRO:HD2	1.86	0.58
13:AN:63:CYS:SG	13:AN:66:THR:HG22	2.44	0.58
23:BB:651:G:P	50:B3:16:THR:HB	2.44	0.58
23:BB:1722:A:H61	23:BB:1738:G:H1'	1.69	0.58
23:BB:18:U:H2'	23:BB:19:A:C8	2.38	0.58
23:BB:2105:U:H2'	23:BB:2106:U:C6	2.39	0.58
23:BB:2378:A:H2'	23:BB:2379:G:H5'	1.85	0.58
23:BB:2415:G:H2'	23:BB:2416:C:C6	2.39	0.58
23:BB:2572:A:C8	26:BD:150:GLN:O	2.57	0.58
23:BB:2783:U:H2'	23:BB:2784:U:H6	1.68	0.58
23:BB:458:G:N2	23:BB:469:G:H2'	2.19	0.58
23:BB:705:A:N6	23:BB:726:G:O2'	2.37	0.58
25:BC:63:ILE:O	25:BC:64:VAL:HG13	2.04	0.58
27:BE:52:VAL:HG13	27:BE:53:THR:N	2.18	0.58
28:BF:132:ARG:HG3	28:BF:132:ARG:HH11	1.68	0.58
23:BB:2530:A:N6	29:BG:155:PRO:HG3	2.19	0.58
29:BG:49:LEU:HD12	29:BG:49:LEU:N	2.19	0.58
30:BH:3:VAL:HG13	30:BH:21:VAL:HG11	1.86	0.58
32:BK:7:MET:HE2	32:BK:20:MET:HA	1.86	0.58
35:BN:116:VAL:HG13	35:BN:117:ASP:N	2.19	0.58
41:BT:58:VAL:HG12	41:BT:59:ASN:H	1.68	0.58
46:BZ:25:ARG:HH11	46:BZ:25:ARG:CB	2.11	0.58
1:CA:1200:C:C3'	1:CA:1201:A:H5'	2.33	0.58
1:CA:398:U:H2'	1:CA:399:G:C8	2.39	0.58
1:CA:687:A:C2	1:CA:704:A:C5	2.91	0.58
1:CA:865:A:H2'	1:CA:866:C:H6	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:129:PHE:CE2	2:CC:156:LEU:HD13	2.38	0.58
6:CG:74:VAL:HG21	6:CG:143:MET:HG2	1.86	0.58
7:CH:77:VAL:HG23	7:CH:126:CYS:HA	1.85	0.58
19:CT:43:LYS:HB3	19:CT:86:ALA:HB3	1.84	0.58
23:DB:1439:A:N7	23:DB:1440:U:N1	2.51	0.58
23:DB:2415:G:H2'	23:DB:2416:C:C6	2.39	0.58
23:DB:267:C:H2'	23:DB:268:C:H6	1.69	0.58
23:DB:2820:A:H4'	35:DN:4:ARG:HG3	1.84	0.58
23:DB:331:C:O2'	23:DB:332:A:H5'	2.04	0.58
23:DB:499:U:H2'	23:DB:500:G:O4'	2.04	0.58
23:DB:634:C:H2'	23:DB:635:C:H6	1.68	0.58
23:DB:936:A:H2'	23:DB:937:C:H6	1.68	0.58
31:DJ:7:LYS:HD2	31:DJ:45:THR:OG1	2.04	0.58
35:DN:11:ASN:C	35:DN:12:ARG:HD3	2.24	0.58
36:DO:75:GLY:HA3	36:DO:109:ALA:HB3	1.86	0.58
23:DB:2257:U:H5'	43:DW:5:ALA:CB	2.34	0.58
1:AA:1060:U:H5''	9:AJ:53:ILE:CG2	2.34	0.58
1:AA:1434:A:H2'	1:AA:1435:G:C8	2.39	0.58
1:AA:501:C:H1'	1:AA:549:C:H1'	1.85	0.58
1:AA:552:U:H2'	1:AA:553:A:H8	1.69	0.58
1:AA:8:A:H1'	4:AE:107:GLY:HA2	1.86	0.58
20:AB:163:ILE:O	20:AB:185:ILE:HG13	2.04	0.58
4:AE:85:LYS:HE2	4:AE:92:ARG:HH11	1.68	0.58
8:AI:21:LYS:O	8:AI:60:LEU:HB2	2.03	0.58
9:AJ:31:ARG:NH1	9:AJ:31:ARG:HG2	2.19	0.58
15:AP:41:PRO:O	15:AP:42:ILE:HD13	2.03	0.58
23:BB:11:C:H2'	23:BB:12:U:H5'	1.85	0.58
23:BB:660:C:H2'	23:BB:661:A:C8	2.39	0.58
26:BD:8:LYS:HD2	37:BP:5:LYS:HD2	1.85	0.58
52:BI:11:GLN:O	52:BI:11:GLN:HG3	2.03	0.58
32:BK:108:ARG:HB2	32:BK:116:ILE:HD13	1.84	0.58
35:BN:67:PHE:HD2	35:BN:73:ASN:HD21	1.52	0.58
36:BO:9:ARG:HA	36:BO:12:THR:HG23	1.85	0.58
41:BT:11:LEU:HB2	41:BT:32:LEU:HD23	1.85	0.58
41:BT:39:THR:O	41:BT:43:ILE:HG12	2.03	0.58
42:BU:40:LEU:HD13	42:BU:41:VAL:H	1.69	0.58
44:BX:22:LEU:HA	44:BX:47:ARG:HH22	1.68	0.58
44:BX:2:LYS:HD2	44:BX:2:LYS:H	1.69	0.58
1:CA:1036:A:H2'	1:CA:1037:C:H6	1.68	0.58
1:CA:1387:G:H2'	1:CA:1388:C:C6	2.38	0.58
1:CA:677:U:H2'	1:CA:678:U:H6	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:801:U:H2'	1:CA:802:A:H8	1.68	0.58
6:CG:12:LEU:HD13	6:CG:13:PRO:HD2	1.85	0.58
14:CO:88:ARG:NH2	23:DB:714:U:H3'	2.19	0.58
49:D2:16:HIS:NE2	49:D2:44:VAL:HA	2.19	0.58
50:D3:49:VAL:HG13	50:D3:51:LYS:H	1.69	0.58
23:DB:1537:G:H3'	23:DB:1537:G:N3	2.19	0.58
23:DB:1847:A:H4'	23:DB:1848:A:C8	2.39	0.58
23:DB:225:C:H2'	23:DB:226:A:O4'	2.03	0.58
23:DB:1818:U:H2'	25:DC:152:GLN:O	2.04	0.58
25:DC:173:LEU:HD12	25:DC:183:VAL:HG11	1.86	0.58
26:DD:116:LYS:HB2	26:DD:165:MET:HB2	1.86	0.58
26:DD:23:PRO:HB3	26:DD:188:LEU:HB3	1.84	0.58
27:DE:172:ALA:O	27:DE:173:THR:HG22	2.04	0.58
27:DE:4:VAL:HG13	27:DE:5:LEU:N	2.14	0.58
29:DG:5:LYS:HD3	29:DG:68:ARG:CB	2.33	0.58
31:DJ:25:LEU:HD12	31:DJ:62:VAL:HA	1.86	0.58
31:DJ:81:ILE:HG23	31:DJ:82:GLY:N	2.12	0.58
23:DB:2839:G:C4'	35:DN:49:GLU:HG2	2.32	0.58
44:DX:49:ASP:HA	44:DX:52:ARG:NH1	2.19	0.58
5:AF:40:GLU:HG3	5:AF:42:TRP:NE1	2.19	0.57
1:AA:537:G:H5''	11:AL:109:ARG:NH1	2.19	0.57
16:AQ:23:ALA:HB1	16:AQ:40:THR:HG23	1.84	0.57
22:BA:32:U:H1'	22:BA:52:A:N7	2.20	0.57
23:BB:1163:G:O2'	23:BB:1164:C:H5'	2.04	0.57
23:BB:1203:U:H3'	23:BB:1204:A:C5'	2.33	0.57
23:BB:1439:A:N7	23:BB:1440:U:N1	2.51	0.57
23:BB:1442:U:H2'	23:BB:1443:U:H6	1.69	0.57
23:BB:1515:A:H4'	23:BB:1556:C:O2'	2.04	0.57
23:BB:1946:U:H2'	23:BB:1947:C:H6	1.69	0.57
23:BB:705:A:N6	23:BB:726:G:H1'	2.19	0.57
23:BB:718:A:H5''	23:BB:719:C:H5	1.69	0.57
23:BB:1818:U:H5''	25:BC:155:ARG:CG	2.34	0.57
27:BE:150:THR:HA	27:BE:187:VAL:HG23	1.85	0.57
27:BE:43:THR:HG23	27:BE:44:ARG:NE	2.19	0.57
12:AM:67:ASP:CG	28:BF:111:ARG:HB3	2.25	0.57
29:BG:20:GLY:HA2	29:BG:41:GLU:HG2	1.86	0.57
30:BH:11:ASN:ND2	30:BH:12:LEU:H	2.01	0.57
31:BJ:10:THR:HB	31:BJ:13:ARG:HH12	1.68	0.57
33:BL:82:LEU:HD13	33:BL:83:ALA:N	2.19	0.57
34:BM:64:TRP:O	34:BM:102:LEU:HB2	2.03	0.57
34:BM:71:LYS:HD2	34:BM:71:LYS:C	2.24	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:12:VAL:HG11	42:BU:18:LYS:H	1.67	0.57
42:BU:23:LYS:HZ3	42:BU:24:VAL:CG2	2.17	0.57
46:BZ:63:ARG:HA	46:BZ:63:ARG:CZ	2.34	0.57
1:CA:1072:G:H2'	1:CA:1073:U:H6	1.69	0.57
1:CA:1096:C:H2'	1:CA:1097:C:H6	1.67	0.57
1:CA:335:C:H2'	1:CA:336:A:C8	2.39	0.57
20:CB:202:ASN:ND2	20:CB:203:ASP:N	2.51	0.57
20:CB:57:ASN:ND2	20:CB:223:GLY:HA2	2.19	0.57
13:CN:40:ARG:HH11	13:CN:40:ARG:HG3	1.69	0.57
15:CP:1:MET:HB3	15:CP:24:SER:OG	2.04	0.57
21:CU:3:ILE:HG12	21:CU:19:LYS:HB3	1.85	0.57
23:DB:125:A:H4'	49:D2:13:ASN:HD21	1.69	0.57
49:D2:46:LYS:O	49:D2:46:LYS:HE3	2.04	0.57
50:D3:2:LYS:HB2	50:D3:2:LYS:NZ	2.19	0.57
23:DB:1552:A:C2'	23:DB:1553:A:H5'	2.33	0.57
23:DB:2072:C:H2'	23:DB:2073:C:H6	1.69	0.57
23:DB:2553:G:H2'	23:DB:2554:U:C4'	2.34	0.57
23:DB:2567:G:H2'	23:DB:2568:U:C6	2.39	0.57
23:DB:2650:U:H2'	23:DB:2651:C:H6	1.69	0.57
25:DC:53:ILE:HG21	25:DC:218:THR:OG1	2.04	0.57
27:DE:141:MET:HB3	27:DE:185:LYS:HZ1	1.68	0.57
29:DG:85:LYS:HB3	29:DG:131:VAL:HA	1.85	0.57
32:DK:105:ARG:H	32:DK:105:ARG:HD3	1.68	0.57
36:DO:27:VAL:HG23	36:DO:28:VAL:H	1.67	0.57
36:DO:30:ARG:N	36:DO:30:ARG:HD3	2.19	0.57
37:DP:27:VAL:HG23	37:DP:86:LYS:HB3	1.85	0.57
38:DQ:92:LYS:C	38:DQ:94:LEU:H	2.07	0.57
40:DS:23:LEU:HB2	47:D0:21:LEU:CD1	2.30	0.57
44:DX:1:MET:HB2	44:DX:6:LEU:HG	1.86	0.57
1:AA:1027:C:H2'	1:AA:1028:C:H6	1.68	0.57
1:AA:1039:G:H2'	1:AA:1040:U:C6	2.39	0.57
1:AA:1491:G:H5''	1:AA:1492:A:OP2	2.04	0.57
20:AB:67:LEU:HD22	20:AB:67:LEU:N	2.19	0.57
2:AC:59:PRO:HG2	2:AC:60:ALA:H	1.69	0.57
23:BB:1123:C:H2'	23:BB:1124:G:H8	1.69	0.57
23:BB:1275:A:C2'	23:BB:1276:A:H5'	2.34	0.57
23:BB:208:C:H2'	23:BB:209:C:H6	1.69	0.57
23:BB:2247:A:H2'	23:BB:2248:C:C6	2.39	0.57
23:BB:2339:C:H2'	23:BB:2340:A:H8	1.69	0.57
23:BB:2553:G:H2'	23:BB:2554:U:C4'	2.34	0.57
23:BB:582:A:H2'	23:BB:583:G:H8	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:BE:118:LEU:HD23	27:BE:118:LEU:N	2.17	0.57
27:BE:121:VAL:HG12	27:BE:122:GLU:N	2.19	0.57
28:BF:64:PRO:HA	28:BF:88:VAL:HG22	1.86	0.57
30:BH:134:VAL:HG13	30:BH:135:HIS:N	2.14	0.57
35:BN:41:ALA:HA	35:BN:113:ILE:HD11	1.87	0.57
35:BN:73:ASN:HA	35:BN:76:VAL:HG13	1.86	0.57
45:BY:4:ILE:HB	45:BY:39:ASP:HB3	1.86	0.57
1:CA:1496:C:H4'	23:DB:1920:C:O2'	2.04	0.57
1:CA:546:A:P	3:CD:68:GLU:HB3	2.44	0.57
7:CH:93:LYS:HG2	7:CH:96:ALA:HA	1.84	0.57
9:CJ:36:VAL:HG13	9:CJ:76:ILE:CD1	2.34	0.57
1:CA:688:G:H5'	10:CK:48:GLY:HA2	1.85	0.57
10:CK:69:CYS:C	10:CK:73:VAL:HG22	2.24	0.57
16:CQ:26:ARG:NH2	16:CQ:39:ARG:HG2	2.19	0.57
51:D4:11:CYS:HB2	51:D4:14:CYS:CB	2.28	0.57
51:D4:15:LYS:HZ2	51:D4:22:VAL:HG12	1.69	0.57
23:DB:1275:A:C2'	23:DB:1276:A:H5'	2.33	0.57
23:DB:135:U:H2'	23:DB:136:G:C8	2.38	0.57
23:DB:155:A:H2'	23:DB:156:A:H8	1.69	0.57
23:DB:1709:U:H2'	23:DB:1710:G:H8	1.68	0.57
23:DB:1739:A:H2'	23:DB:1740:G:C8	2.40	0.57
23:DB:179:C:H2'	23:DB:180:G:O4'	2.03	0.57
23:DB:526:A:N6	23:DB:2626:C:C4'	2.66	0.57
23:DB:660:C:H2'	23:DB:661:A:H8	1.68	0.57
23:DB:921:C:H2'	23:DB:922:C:H6	1.68	0.57
23:DB:962:G:H21	23:DB:2250:G:H1	1.52	0.57
23:DB:987:C:H2'	23:DB:988:A:O4'	2.04	0.57
26:DD:204:LYS:HE2	26:DD:204:LYS:HA	1.84	0.57
23:DB:536:G:H5''	38:DQ:52:ARG:HH22	1.69	0.57
23:DB:1198:U:H4'	38:DQ:8:ILE:HD11	1.87	0.57
23:DB:2256:G:O2'	43:DW:5:ALA:HB1	2.03	0.57
45:DY:37:ARG:CZ	45:DY:37:ARG:HA	2.34	0.57
46:DZ:47:LYS:HB2	46:DZ:51:VAL:CG1	2.34	0.57
1:AA:335:C:H2'	1:AA:336:A:C8	2.39	0.57
2:AC:77:GLY:HA3	2:AC:81:GLU:HB3	1.87	0.57
8:AI:10:ARG:HB2	8:AI:14:SER:O	2.04	0.57
9:AJ:15:HIS:HD2	9:AJ:18:ILE:HG22	1.69	0.57
47:B0:52:LYS:HZ3	47:B0:52:LYS:HA	1.68	0.57
50:B3:24:LYS:NZ	50:B3:29:ARG:HH22	2.02	0.57
23:BB:1560:G:H2'	23:BB:1561:C:H6	1.70	0.57
23:BB:2352:A:H2'	23:BB:2353:G:O4'	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:346:A:H2'	23:BB:347:A:O4'	2.04	0.57
23:BB:404:A:H4'	23:BB:405:U:H5'	1.87	0.57
26:BD:2:ILE:HG12	26:BD:203:VAL:O	2.04	0.57
27:BE:120:VAL:H	27:BE:189:THR:CG2	2.17	0.57
27:BE:128:ALA:CB	27:BE:129:PRO:CD	2.81	0.57
37:BP:25:VAL:HG23	37:BP:49:ILE:HD11	1.84	0.57
43:BW:49:ASN:HB2	43:BW:53:GLY:CA	2.34	0.57
1:CA:114:U:O2'	1:CA:115:G:H5'	2.04	0.57
1:CA:441:A:H61	1:CA:493:A:N6	2.02	0.57
1:CA:552:U:H2'	1:CA:553:A:H8	1.70	0.57
1:CA:1057:G:O3'	2:CC:196:GLY:HA3	2.04	0.57
3:CD:115:GLN:NE2	3:CD:153:ARG:HH22	2.03	0.57
4:CE:72:ASN:N	4:CE:72:ASN:HD22	2.01	0.57
10:CK:12:ARG:HB3	10:CK:13:LYS:NZ	2.19	0.57
18:CS:22:VAL:HG22	18:CS:42:ASN:OD1	2.04	0.57
49:D2:12:ARG:HH21	49:D2:16:HIS:HB2	1.69	0.57
23:DB:1082:U:C2	23:DB:1086:A:C6	2.92	0.57
23:DB:182:A:H2'	23:DB:183:C:C6	2.39	0.57
23:DB:230:G:H2'	23:DB:231:A:C8	2.39	0.57
23:DB:2590:A:H2'	23:DB:2591:C:C6	2.40	0.57
23:DB:301:G:H3'	23:DB:335:C:OP2	2.04	0.57
23:DB:582:A:H2'	23:DB:583:G:H8	1.68	0.57
23:DB:72:U:O2'	23:DB:73:A:H5'	2.04	0.57
26:DD:174:SER:C	26:DD:175:LEU:HD12	2.25	0.57
26:DD:22:ILE:HG23	26:DD:191:GLY:N	2.19	0.57
30:DH:26:ALA:HB3	30:DH:31:VAL:CG2	2.33	0.57
36:DO:90:VAL:HG13	36:DO:116:GLN:H	1.68	0.57
37:DP:25:VAL:HG12	37:DP:27:VAL:N	2.18	0.57
32:DK:76:VAL:HB	37:DP:74:GLN:NE2	2.17	0.57
41:DT:58:VAL:HG12	41:DT:59:ASN:H	1.69	0.57
42:DU:23:LYS:C	42:DU:25:LYS:HD2	2.24	0.57
42:DU:38:ILE:HG13	42:DU:62:ALA:HB1	1.87	0.57
44:DX:55:THR:HG22	44:DX:56:LEU:H	1.69	0.57
45:DY:18:LYS:CD	45:DY:18:LYS:H	2.11	0.57
46:DZ:55:GLY:HA2	46:DZ:59:ARG:HD2	1.86	0.57
1:AA:1260:G:H4'	1:AA:1284:C:H5'	1.86	0.57
1:AA:1508:A:H2'	1:AA:1509:C:H6	1.69	0.57
1:AA:450:G:N7	1:AA:481:G:O6	2.37	0.57
1:AA:950:U:H2'	1:AA:951:G:C8	2.39	0.57
3:AD:55:ARG:HH21	3:AD:58:GLN:NE2	2.02	0.57
6:AG:14:ASP:OD2	6:AG:22:LEU:HD12	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:187:G:H4'	19:AT:79:THR:HG21	1.86	0.57
50:B3:24:LYS:HZ3	50:B3:29:ARG:NH2	2.02	0.57
22:BA:50:A:H5''	36:BO:68:LYS:CD	2.33	0.57
23:BB:727:A:OP1	23:BB:1431:A:O2'	2.21	0.57
23:BB:15:G:O2'	23:BB:16:C:H5'	2.05	0.57
23:BB:2307:G:H4'	23:BB:2311:A:H61	1.68	0.57
26:BD:2:ILE:HB	26:BD:205:PRO:HD3	1.85	0.57
29:BG:84:LYS:HZ2	29:BG:85:LYS:H	1.50	0.57
31:BJ:25:LEU:N	31:BJ:25:LEU:HD13	2.19	0.57
34:BM:108:VAL:N	34:BM:109:PRO:HD3	2.19	0.57
35:BN:10:LEU:H	35:BN:17:ARG:NH1	2.02	0.57
35:BN:29:VAL:HG21	35:BN:75:ILE:HD11	1.86	0.57
36:BO:40:ILE:H	36:BO:40:ILE:HD13	1.69	0.57
40:BS:18:ARG:NH1	40:BS:25:ARG:HH22	2.02	0.57
41:BT:10:VAL:O	41:BT:34:VAL:HG13	2.05	0.57
24:BV:80:HIS:N	24:BV:87:GLN:HE22	2.02	0.57
1:CA:1105:A:H2'	1:CA:1106:G:H8	1.69	0.57
1:CA:1461:G:H2'	1:CA:1462:C:C6	2.39	0.57
1:CA:265:G:H2'	1:CA:267:C:H5	1.69	0.57
1:CA:383:A:H2'	1:CA:384:G:O4'	2.05	0.57
1:CA:411:A:C4	1:CA:413:G:H1'	2.40	0.57
4:CE:73:VAL:HG21	4:CE:143:LEU:O	2.04	0.57
14:CO:53:ARG:HG2	14:CO:53:ARG:HH11	1.68	0.57
51:D4:16:ILE:HG23	51:D4:18:LYS:H	1.69	0.57
23:DB:1932:A:H2'	23:DB:1933:G:O4'	2.04	0.57
23:DB:196:A:N3	23:DB:196:A:H2'	2.19	0.57
23:DB:2053:G:H5'	26:DD:149:ASN:O	2.04	0.57
23:DB:2283:C:H5''	23:DB:2389:G:O2'	2.04	0.57
23:DB:2376:A:N3	36:DO:111:ARG:NH2	2.52	0.57
23:DB:2617:U:C2'	23:DB:2618:G:H5'	2.35	0.57
23:DB:2819:G:H2'	23:DB:2821:A:N7	2.19	0.57
23:DB:849:A:H2'	23:DB:850:U:C6	2.40	0.57
23:DB:877:A:H2	23:DB:900:A:N7	2.02	0.57
30:DH:113:SER:HB2	30:DH:132:PHE:HZ	1.69	0.57
52:DI:27:LEU:HD23	52:DI:27:LEU:N	2.15	0.57
33:DL:30:THR:HG21	33:DL:38:GLN:NE2	2.20	0.57
23:DB:536:G:H5''	38:DQ:52:ARG:NH2	2.20	0.57
39:DR:80:ARG:CZ	39:DR:85:LYS:HB3	2.34	0.57
1:AA:1366:C:H2'	1:AA:1367:C:C6	2.39	0.57
13:AN:66:THR:HG23	13:AN:67:GLY:H	1.69	0.57
1:AA:1320:C:H41	18:AS:36:ARG:CG	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:B1:15:GLY:N	48:B1:47:ILE:HD12	2.19	0.57
23:BB:1438:U:N3	23:BB:1552:A:N6	2.53	0.57
23:BB:1505:A:H2'	23:BB:1506:U:C6	2.39	0.57
23:BB:1537:G:H2'	23:BB:1538:G:C4'	2.35	0.57
23:BB:1878:G:H2'	23:BB:1879:C:H6	1.69	0.57
23:BB:264:C:O2'	23:BB:265:A:H5''	2.05	0.57
23:BB:285:G:O2'	23:BB:286:U:H5'	2.05	0.57
23:BB:716:A:H2'	23:BB:717:C:O4'	2.03	0.57
23:BB:918:A:C2'	23:BB:919:U:H5'	2.34	0.57
25:BC:76:VAL:O	25:BC:93:VAL:HA	2.04	0.57
27:BE:108:ILE:HG21	27:BE:116:ASP:OD2	2.04	0.57
27:BE:116:ASP:O	27:BE:117:ARG:HB3	2.05	0.57
27:BE:48:THR:HG22	27:BE:49:ARG:N	2.14	0.57
28:BF:177:ARG:HE	28:BF:178:LYS:CA	2.17	0.57
31:BJ:25:LEU:HG	31:BJ:63:ALA:H	1.69	0.57
34:BM:11:LYS:O	34:BM:12:MET:HE2	2.03	0.57
34:BM:34:LYS:HD3	34:BM:98:PRO:O	2.04	0.57
36:BO:17:LYS:HE3	36:BO:43:ASN:HB2	1.86	0.57
41:BT:3:ARG:HD3	41:BT:3:ARG:C	2.23	0.57
42:BU:78:LYS:HG3	42:BU:96:LYS:HE3	1.85	0.57
24:BV:84:PRO:HG3	34:BM:127:LYS:NZ	2.19	0.57
43:BW:49:ASN:HD22	43:BW:53:GLY:HA2	1.68	0.57
1:CA:238:A:C2'	1:CA:239:U:H5''	2.31	0.57
1:CA:972:C:OP1	9:CJ:59:LYS:HD2	2.04	0.57
20:CB:169:HIS:CE1	20:CB:173:LYS:HB2	2.40	0.57
3:CD:58:GLN:O	3:CD:62:ARG:HG2	2.04	0.57
9:CJ:37:ARG:HG2	9:CJ:37:ARG:HH11	1.68	0.57
9:CJ:36:VAL:HA	9:CJ:76:ILE:HG23	1.87	0.57
12:CM:48:SER:H	12:CM:51:GLN:NE2	2.03	0.57
21:CU:41:THR:CA	21:CU:45:LYS:HD3	2.32	0.57
51:D4:26:ILE:O	51:D4:27:CYS:CB	2.52	0.57
22:DA:59:A:H2'	22:DA:60:C:O4'	2.05	0.57
23:DB:1637:A:H2'	23:DB:1638:C:C6	2.39	0.57
23:DB:1791:A:H5''	25:DC:211:ARG:HE	1.70	0.57
23:DB:2085:U:O2'	23:DB:2086:U:H5'	2.05	0.57
23:DB:2329:U:H2'	23:DB:2330:G:H8	1.69	0.57
23:DB:2869:G:H2'	23:DB:2870:C:C6	2.40	0.57
52:DI:37:PHE:CZ	52:DI:58:ILE:HD11	2.39	0.57
52:DI:45:THR:HA	52:DI:48:ILE:CG2	2.33	0.57
33:DL:118:THR:O	33:DL:119:PRO:C	2.39	0.57
33:DL:81:ASP:HA	33:DL:84:LYS:HB3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:5:LYS:HA	34:DM:68:PHE:CE1	2.40	0.57
36:DO:18:LEU:HD21	36:DO:43:ASN:ND2	2.19	0.57
37:DP:47:ILE:CG2	37:DP:48:ALA:H	2.03	0.57
42:DU:12:VAL:HG11	42:DU:17:ASP:HB3	1.87	0.57
24:DV:4:ILE:N	24:DV:4:ILE:HD12	2.19	0.57
43:DW:55:ASP:CG	43:DW:56:HIS:H	2.08	0.57
1:AA:1074:G:H5'	20:AB:104:LYS:NZ	2.19	0.57
1:AA:1283:U:H2'	1:AA:1284:C:C6	2.40	0.57
1:AA:202:G:H2'	1:AA:203:G:H8	1.70	0.57
1:AA:411:A:C4	1:AA:413:G:H1'	2.40	0.57
1:AA:71:A:N1	1:AA:99:C:H1'	2.19	0.57
4:AE:81:GLN:HE22	4:AE:82:HIS:CE1	2.23	0.57
6:AG:14:ASP:H	6:AG:19:SER:H	1.51	0.57
16:AQ:7:LEU:HD13	16:AQ:24:ILE:HG12	1.87	0.57
18:AS:11:ASP:H	18:AS:37:SER:HB3	1.68	0.57
21:AU:39:LYS:N	21:AU:40:PRO:CD	2.66	0.57
51:B4:22:VAL:HG11	51:B4:36:ARG:NH1	2.20	0.57
23:BB:1180:U:O2'	23:BB:1181:U:H5'	2.05	0.57
23:BB:1637:A:H2'	23:BB:1638:C:C6	2.40	0.57
23:BB:2083:G:H2'	23:BB:2084:C:H6	1.69	0.57
23:BB:548:G:H5''	23:BB:549:G:O4'	2.03	0.57
23:BB:871:U:H2'	23:BB:872:U:H6	1.68	0.57
27:BE:112:LEU:HB3	27:BE:116:ASP:H	1.69	0.57
27:BE:24:ASN:O	27:BE:24:ASN:CG	2.43	0.57
33:BL:78:ARG:NH1	33:BL:110:VAL:HG11	2.20	0.57
33:BL:79:LEU:HB3	33:BL:114:GLY:N	2.12	0.57
34:BM:96:ILE:HD13	34:BM:97:GLN:CG	2.34	0.57
37:BP:62:LYS:HD3	37:BP:72:VAL:HG21	1.87	0.57
38:BQ:51:GLN:O	38:BQ:54:ARG:HG2	2.05	0.57
22:DA:113:C:O2'	36:DO:47:VAL:HA	2.04	0.57
23:DB:1309:G:C4'	49:D2:7:PRO:HB2	2.34	0.57
23:DB:1441:G:H2'	23:DB:1442:U:C6	2.40	0.57
23:DB:1464:G:H2'	23:DB:1465:G:H8	1.68	0.57
23:DB:1505:A:H2'	23:DB:1506:U:C6	2.39	0.57
23:DB:2105:U:H2'	23:DB:2106:U:H6	1.68	0.57
23:DB:2480:C:O2'	23:DB:2481:G:H5'	2.03	0.57
23:DB:2722:G:O2'	35:DN:4:ARG:NE	2.38	0.57
23:DB:453:A:H4'	23:DB:472:A:H61	1.69	0.57
25:DC:156:SER:O	25:DC:195:GLY:HA3	2.05	0.57
25:DC:179:GLU:CG	25:DC:266:ILE:HA	2.34	0.57
26:DD:122:VAL:CA	26:DD:128:ARG:HG3	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:133:LEU:HD22	27:DE:136:GLN:HG3	1.87	0.57
28:DF:120:SER:HB3	28:DF:127:TYR:CE1	2.40	0.57
29:DG:124:CYS:HB3	29:DG:130:ILE:HD13	1.87	0.57
30:DH:6:LEU:HB2	30:DH:35:LYS:HB2	1.85	0.57
30:DH:96:THR:HG23	30:DH:97:ARG:H	1.68	0.57
52:DI:100:ILE:O	52:DI:139:VAL:HA	2.04	0.57
32:DK:70:ARG:CD	32:DK:76:VAL:HG22	2.33	0.57
33:DL:133:ALA:O	33:DL:136:GLU:HB2	2.04	0.57
34:DM:33:LEU:HD22	34:DM:124:LEU:HD22	1.85	0.57
37:DP:1:SER:HA	37:DP:4:ILE:HB	1.86	0.57
37:DP:27:VAL:HA	37:DP:86:LYS:CE	2.32	0.57
38:DQ:65:ASN:HB2	38:DQ:75:TYR:HB2	1.87	0.57
38:DQ:111:LYS:NZ	39:DR:52:PRO:HA	2.20	0.57
40:DS:87:PRO:O	40:DS:93:ALA:HA	2.05	0.57
36:DO:15:ARG:HH22	43:DW:76:ARG:CD	2.16	0.57
1:AA:114:U:O2'	1:AA:115:G:H5'	2.04	0.57
1:AA:1376:U:H2'	1:AA:1377:A:H8	1.69	0.57
1:AA:413:G:H4'	1:AA:414:A:H5''	1.85	0.57
1:AA:85:U:H4'	1:AA:86:G:H4'	1.85	0.57
1:AA:973:G:H3'	1:AA:974:A:H5''	1.87	0.57
20:AB:102:ASN:O	20:AB:106:VAL:HG23	2.04	0.57
2:AC:21:TRP:HB3	2:AC:58:ARG:HB2	1.86	0.57
3:AD:25:ARG:HH11	3:AD:30:LYS:HE3	1.69	0.57
1:AA:1179:A:H4'	8:AI:104:THR:HA	1.86	0.57
8:AI:107:ALA:O	8:AI:109:GLN:HG2	2.05	0.57
10:AK:69:CYS:C	10:AK:71:ASP:H	2.06	0.57
10:AK:95:THR:HG23	10:AK:96:ILE:N	2.19	0.57
11:AL:34:THR:HB	11:AL:53:ARG:HB2	1.87	0.57
21:AU:11:PHE:O	21:AU:11:PHE:HD1	1.86	0.57
51:B4:33:HIS:HD2	51:B4:33:HIS:H	1.51	0.57
23:BB:1220:G:H2'	23:BB:1221:C:C6	2.40	0.57
23:BB:2650:U:H2'	23:BB:2651:C:H6	1.69	0.57
23:BB:2819:G:H2'	23:BB:2821:A:N7	2.19	0.57
23:BB:2872:A:O2'	23:BB:2873:A:H5''	2.05	0.57
23:BB:2882:A:H2'	23:BB:2883:A:H5''	1.86	0.57
23:BB:519:U:H5''	40:BS:18:ARG:NH2	2.20	0.57
23:BB:632:A:H2'	23:BB:633:A:C8	2.40	0.57
28:BF:68:LYS:HA	28:BF:83:PRO:HA	1.86	0.57
32:BK:53:LYS:HE2	32:BK:56:ASP:OD1	2.04	0.57
40:BS:6:LYS:HA	40:BS:104:THR:CA	2.28	0.57
41:BT:64:LYS:HB3	41:BT:77:ARG:HD3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BV:44:HIS:NE2	24:BV:85:LYS:HB2	2.19	0.57
43:BW:36:ILE:O	43:BW:36:ILE:HG22	2.05	0.57
46:BZ:50:ASP:HA	46:BZ:53:THR:HG22	1.86	0.57
1:CA:1175:G:O2'	1:CA:1176:A:H5'	2.04	0.57
1:CA:384:G:H2'	1:CA:385:C:H6	1.68	0.57
4:CE:142:GLY:HA2	4:CE:145:ASN:HD22	1.70	0.57
5:CF:38:ARG:NH2	5:CF:96:VAL:HB	2.20	0.57
7:CH:101:ALA:O	7:CH:103:VAL:HG23	2.04	0.57
13:CN:40:ARG:HG3	18:CS:6:LYS:O	2.05	0.57
15:CP:54:LEU:HD22	15:CP:80:LYS:HZ2	1.70	0.57
22:DA:112:G:O2'	22:DA:113:C:H5'	2.04	0.57
23:DB:1381:G:O2'	23:DB:1382:G:H5'	2.05	0.57
23:DB:1400:U:H2'	23:DB:1401:G:H8	1.70	0.57
23:DB:1722:A:H61	23:DB:1738:G:H1'	1.69	0.57
23:DB:1729:U:H2'	23:DB:1730:C:H4'	1.87	0.57
23:DB:1842:G:H2'	23:DB:1843:C:H6	1.70	0.57
23:DB:2352:A:H2'	23:DB:2353:G:O4'	2.04	0.57
23:DB:2455:G:H2'	23:DB:2456:C:H6	1.70	0.57
23:DB:558:U:P	31:DJ:113:PRO:HB2	2.45	0.57
25:DC:53:ILE:HD13	25:DC:218:THR:HG23	1.87	0.57
27:DE:109:LEU:HG	27:DE:117:ARG:CD	2.34	0.57
29:DG:163:TYR:HB2	29:DG:166:GLU:HG3	1.85	0.57
31:DJ:89:PHE:HD1	31:DJ:92:MET:HG3	1.68	0.57
33:DL:109:LYS:HA	33:DL:127:VAL:H	1.69	0.57
35:DN:35:LYS:HE3	35:DN:110:MET:HB3	1.86	0.57
36:DO:83:LEU:HA	36:DO:87:ILE:HD12	1.87	0.57
41:DT:3:ARG:HB3	41:DT:5:GLU:OE2	2.04	0.57
24:DV:82:TYR:HE1	24:DV:83:LYS:HE3	1.70	0.57
43:DW:42:THR:HG23	43:DW:66:VAL:N	2.18	0.57
44:DX:17:GLU:HA	44:DX:21:LEU:HB2	1.85	0.57
45:DY:10:ARG:O	45:DY:11:SER:HB3	2.04	0.57
1:AA:1081:A:O2'	1:AA:1082:A:H5'	2.05	0.57
1:AA:537:G:H5''	11:AL:109:ARG:HH11	1.69	0.57
1:AA:1101:A:H61	20:AB:173:LYS:HD2	1.70	0.57
5:AF:3:HIS:HB2	5:AF:92:THR:CB	2.35	0.57
8:AI:113:LYS:HA	8:AI:120:ALA:HB2	1.87	0.57
9:AJ:42:LEU:HB2	9:AJ:71:LEU:HD13	1.85	0.57
23:BB:127:A:H5''	23:BB:128:C:C6	2.40	0.57
23:BB:2455:G:H2'	23:BB:2456:C:H6	1.70	0.57
23:BB:643:A:H61	23:BB:2370:G:H1'	1.68	0.57
27:BE:52:VAL:C	27:BE:54:GLY:H	2.08	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:44:ALA:HB1	28:BF:48:LEU:HD12	1.86	0.57
36:BO:17:LYS:HE2	43:BW:77:LYS:CD	2.35	0.57
38:BQ:32:ARG:O	38:BQ:36:GLN:HB2	2.05	0.57
42:BU:11:ILE:HB	42:BU:69:VAL:CB	2.35	0.57
23:BB:2270:A:H4'	43:BW:18:LYS:CB	2.35	0.57
1:CA:1414:U:H2'	1:CA:1415:G:H8	1.69	0.57
51:D4:24:ARG:HH21	51:D4:37:GLN:HB2	1.69	0.57
23:DB:1198:U:H2'	23:DB:1199:U:H6	1.70	0.57
23:DB:1198:U:H2'	23:DB:1199:U:C6	2.39	0.57
23:DB:165:A:H2'	23:DB:166:U:H6	1.69	0.57
23:DB:170:U:H2'	23:DB:171:U:C6	2.39	0.57
23:DB:1946:U:H2'	23:DB:1947:C:H6	1.69	0.57
23:DB:2378:A:H2'	23:DB:2379:G:H5'	1.85	0.57
23:DB:2902:C:O2'	23:DB:2903:U:H4'	2.05	0.57
23:DB:506:G:H4'	23:DB:509:C:O2	2.05	0.57
25:DC:62:ARG:O	25:DC:63:ILE:HB	2.05	0.57
28:DF:36:ASN:HB2	28:DF:87:LYS:HA	1.87	0.57
23:DB:1100:C:OP2	52:DI:2:LYS:HB3	2.05	0.57
52:DI:32:VAL:HG22	52:DI:60:VAL:CG2	2.34	0.57
38:DQ:48:ASP:CA	38:DQ:51:GLN:HE21	2.06	0.57
43:DW:24:ARG:HB3	43:DW:59:PHE:CD2	2.39	0.57
1:AA:212:G:H2'	1:AA:213:G:H8	1.68	0.57
20:AB:37:VAL:HG22	20:AB:38:HIS:H	1.68	0.57
20:AB:66:ILE:C	20:AB:67:LEU:HD13	2.25	0.57
20:AB:71:THR:CG2	20:AB:94:ARG:HH21	2.18	0.57
2:AC:33:ASP:OD2	13:AN:64:ARG:HD3	2.05	0.57
7:AH:29:SER:HB3	7:AH:32:LYS:CG	2.35	0.57
10:AK:126:ARG:HB2	21:AU:33:ARG:CD	2.30	0.57
15:AP:25:ARG:HD3	15:AP:25:ARG:N	2.20	0.57
18:AS:28:LYS:HB2	18:AS:29:PRO:HD2	1.87	0.57
23:BB:1400:U:H2'	23:BB:1401:G:H8	1.70	0.57
23:BB:2213:U:C2'	23:BB:2213:U:O2	2.52	0.57
23:BB:2567:G:H2'	23:BB:2568:U:C6	2.40	0.57
23:BB:2839:G:H2'	23:BB:2840:C:H6	1.70	0.57
23:BB:459:U:C5'	49:B2:41:ARG:HE	2.16	0.57
23:BB:49:A:OP1	23:BB:51:G:H5'	2.05	0.57
25:BC:143:VAL:HG12	25:BC:144:GLU:N	2.14	0.57
25:BC:224:MET:HG3	25:BC:233:GLY:O	2.05	0.57
26:BD:193:VAL:HG13	26:BD:193:VAL:O	2.03	0.57
32:BK:70:ARG:CB	32:BK:76:VAL:HG22	2.31	0.57
33:BL:129:LYS:O	33:BL:129:LYS:HD3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:22:ARG:HG3	35:BN:23:ASN:N	2.17	0.57
37:BP:55:HIS:O	37:BP:57:ALA:N	2.38	0.57
37:BP:62:LYS:HE3	37:BP:74:GLN:NE2	2.19	0.57
38:BQ:50:ARG:HG3	38:BQ:53:LYS:NZ	2.20	0.57
38:BQ:85:ALA:HB1	38:BQ:88:GLU:HB2	1.87	0.57
39:BR:49:ILE:HG22	39:BR:51:VAL:HG23	1.86	0.57
39:BR:58:VAL:HG22	39:BR:59:ILE:N	2.17	0.57
40:BS:1:MET:HB3	40:BS:4:ILE:HG12	1.86	0.57
42:BU:9:GLU:HA	42:BU:23:LYS:HA	1.86	0.57
42:BU:2:ALA:O	42:BU:4:ILE:N	2.37	0.57
42:BU:79:ALA:O	42:BU:96:LYS:HA	2.04	0.57
24:BV:61:LEU:O	24:BV:71:LYS:HA	2.05	0.57
41:BT:12:ARG:NH2	44:BX:29:ARG:HH12	2.01	0.57
44:BX:29:ARG:N	44:BX:34:SER:HB2	2.19	0.57
1:CA:1134:G:C6	1:CA:1135:U:H1'	2.40	0.57
1:CA:908:A:H2'	1:CA:909:A:H8	1.70	0.57
2:CC:115:VAL:O	2:CC:119:ILE:HG22	2.05	0.57
2:CC:143:LEU:HD13	2:CC:143:LEU:H	1.70	0.57
1:CA:619:U:H3	3:CD:130:ASN:HD21	1.53	0.57
8:CI:93:LEU:O	8:CI:97:LEU:HG	2.05	0.57
23:DB:1211:C:H4'	23:DB:1212:G:OP2	2.04	0.57
23:DB:125:A:H4'	23:DB:126:A:OP2	2.05	0.57
23:DB:1537:G:H2'	23:DB:1538:G:C4'	2.35	0.57
23:DB:1791:A:C2	23:DB:1829:A:C4'	2.88	0.57
23:DB:1853:A:N1	23:DB:2087:G:H1'	2.19	0.57
23:DB:2234:G:O2'	23:DB:2235:G:H5'	2.04	0.57
23:DB:544:C:O2'	23:DB:545:U:O5'	2.23	0.57
23:DB:546:U:O2	23:DB:546:U:H5'	2.05	0.57
23:DB:765:C:H2'	23:DB:766:U:C6	2.40	0.57
25:DC:237:ARG:HH11	25:DC:238:ASN:HA	1.69	0.57
28:DF:106:ALA:HB1	28:DF:136:ILE:HG22	1.85	0.57
32:DK:24:VAL:HA	32:DK:39:ILE:HD12	1.86	0.57
32:DK:6:THR:O	32:DK:20:MET:HG3	2.05	0.57
23:DB:962:G:N2	34:DM:81:ARG:HD3	2.20	0.57
38:DQ:69:ARG:HH11	38:DQ:69:ARG:CB	2.16	0.57
45:DY:8:GLN:HB2	45:DY:28:LEU:HG	1.87	0.57
1:AA:265:G:H2'	1:AA:267:C:H5	1.69	0.57
1:AA:266:G:HO2'	1:AA:267:C:H3'	1.70	0.57
1:AA:600:A:OP1	7:AH:88:LYS:HG2	2.04	0.57
1:AA:784:A:H2'	1:AA:785:G:H8	1.70	0.57
1:AA:993:G:H4'	1:AA:994:A:OP2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AB:165:ALA:HB2	20:AB:186:VAL:HG12	1.87	0.57
2:AC:46:LEU:HD23	2:AC:75:VAL:HG13	1.86	0.57
3:AD:99:ASN:CG	3:AD:103:ARG:HH21	2.07	0.57
3:AD:170:LEU:HD12	3:AD:170:LEU:O	2.05	0.57
4:AE:155:LYS:HA	7:AH:65:PHE:CD1	2.40	0.57
8:AI:11:ARG:NH2	8:AI:12:LYS:HE3	2.20	0.57
8:AI:44:ARG:N	8:AI:44:ARG:HH11	2.02	0.57
12:AM:21:ILE:HB	12:AM:24:VAL:HG22	1.87	0.57
13:AN:17:ASP:HA	13:AN:21:ALA:HB2	1.87	0.57
1:AA:1225:A:H4'	18:AS:77:ARG:NH2	2.20	0.57
47:B0:24:VAL:O	47:B0:25:THR:HG23	2.05	0.57
49:B2:30:VAL:HA	49:B2:33:ARG:HD2	1.87	0.57
23:BB:1155:A:OP1	38:BQ:54:ARG:HD2	2.05	0.57
23:BB:1790:C:H4'	25:BC:207:ALA:HB2	1.86	0.57
23:BB:182:A:H2'	23:BB:183:C:H6	1.70	0.57
23:BB:2371:G:O2'	48:B1:44:GLN:HA	2.04	0.57
23:BB:2617:U:C2'	23:BB:2618:G:H5'	2.35	0.57
23:BB:278:A:H2'	23:BB:278:A:N3	2.20	0.57
23:BB:609:A:H2'	23:BB:610:C:O4'	2.05	0.57
23:BB:702:U:O2'	23:BB:703:U:H5'	2.05	0.57
36:BO:43:ASN:HA	36:BO:46:GLU:OE2	2.04	0.57
36:BO:60:GLU:HG3	36:BO:61:GLN:H	1.70	0.57
42:BU:81:ARG:N	42:BU:81:ARG:HD2	2.19	0.57
43:BW:16:GLU:HG2	43:BW:37:VAL:HG22	1.85	0.57
1:CA:51:A:O2'	1:CA:52:C:OP2	2.18	0.57
1:CA:674:G:H2'	1:CA:675:A:C8	2.40	0.57
8:CI:56:MET:SD	8:CI:57:VAL:N	2.78	0.57
9:CJ:52:LEU:HG	9:CJ:62:ARG:HE	1.70	0.57
23:DB:2339:C:H2'	23:DB:2340:A:H8	1.68	0.57
23:DB:632:A:H2'	23:DB:633:A:C8	2.40	0.57
23:DB:796:C:H2'	23:DB:797:G:C8	2.40	0.57
23:DB:997:G:C5'	38:DQ:92:LYS:HG3	2.35	0.57
25:DC:163:ILE:HG12	25:DC:173:LEU:HD21	1.87	0.57
23:DB:2820:A:C6	26:DD:197:THR:HB	2.40	0.57
27:DE:91:ASP:O	27:DE:92:HIS:HB2	2.05	0.57
52:DI:23:VAL:HG12	52:DI:27:LEU:HD21	1.87	0.57
31:DJ:117:ALA:C	31:DJ:121:LYS:HZ3	2.09	0.57
33:DL:55:MET:CB	33:DL:56:PRO:HD3	2.35	0.57
35:DN:73:ASN:HA	35:DN:76:VAL:CG1	2.35	0.57
1:AA:441:A:H61	1:AA:493:A:N6	2.03	0.56
20:AB:202:ASN:C	20:AB:202:ASN:HD22	2.08	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AE:108:GLY:H	4:AE:110:MET:HE1	1.69	0.56
4:AE:92:ARG:HB3	4:AE:127:TYR:HB2	1.87	0.56
6:AG:139:ASP:HA	6:AG:142:ARG:HH12	1.70	0.56
1:AA:1342:C:H5'	8:AI:127:SER:HA	1.86	0.56
8:AI:23:GLY:O	8:AI:24:ASN:HB2	2.04	0.56
9:AJ:52:LEU:HB2	13:AN:80:ARG:HD2	1.87	0.56
16:AQ:56:ASP:HB3	16:AQ:79:GLU:O	2.05	0.56
49:B2:25:LYS:HE3	49:B2:25:LYS:H	1.68	0.56
23:BB:1083:U:H1'	23:BB:1086:A:H61	1.70	0.56
23:BB:1441:G:H2'	23:BB:1442:U:C6	2.40	0.56
23:BB:1464:G:H2'	23:BB:1465:G:H8	1.68	0.56
23:BB:1754:A:OP1	37:BP:95:LYS:HB2	2.04	0.56
23:BB:1771:C:H2'	23:BB:1772:A:H8	1.69	0.56
23:BB:2329:U:H2'	23:BB:2330:G:C8	2.40	0.56
23:BB:2443:C:O2'	23:BB:2444:G:H5'	2.05	0.56
25:BC:21:PRO:C	25:BC:202:ARG:HD2	2.25	0.56
28:BF:108:PRO:HA	28:BF:113:PHE:CD1	2.40	0.56
30:BH:2:GLN:N	30:BH:21:VAL:HG12	2.20	0.56
52:BI:85:ILE:CD1	52:BI:137:LEU:HD21	2.35	0.56
31:BJ:10:THR:HG22	31:BJ:50:THR:N	2.20	0.56
31:BJ:61:LYS:CA	31:BJ:61:LYS:HE2	2.35	0.56
34:BM:101:VAL:O	34:BM:102:LEU:HD12	2.05	0.56
34:BM:41:LEU:HB2	34:BM:93:VAL:CG2	2.35	0.56
24:BV:63:ILE:HD11	24:BV:72:VAL:HG21	1.87	0.56
23:BB:2258:C:OP2	43:BW:3:LYS:HG3	2.05	0.56
43:BW:43:LYS:HG2	43:BW:76:ARG:NH2	2.18	0.56
46:BZ:25:ARG:HG3	46:BZ:26:SER:H	1.70	0.56
1:CA:1011:C:H2'	1:CA:1012:A:C8	2.40	0.56
1:CA:1034:G:H2'	1:CA:1035:A:C8	2.40	0.56
1:CA:1330:U:H2'	1:CA:1331:G:O4'	2.05	0.56
1:CA:784:A:H2'	1:CA:785:G:H8	1.70	0.56
1:CA:940:C:H2'	1:CA:941:G:H8	1.70	0.56
1:CA:1101:A:H61	20:CB:101:THR:HG21	1.70	0.56
2:CC:49:ALA:HB2	2:CC:74:ILE:CG2	2.34	0.56
3:CD:84:ASN:ND2	3:CD:87:GLU:H	2.02	0.56
9:CJ:53:ILE:HG13	13:CN:84:ARG:CZ	2.35	0.56
14:CO:76:ARG:O	14:CO:79:ARG:HB2	2.05	0.56
1:CA:673:A:H1'	17:CR:63:TYR:CD1	2.39	0.56
17:CR:72:ARG:NH1	17:CR:72:ARG:HB2	2.19	0.56
48:D1:19:PHE:CE2	48:D1:41:VAL:HG13	2.39	0.56
23:DB:1022:G:N2	23:DB:1142:A:N1	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1560:G:H2'	23:DB:1561:C:H6	1.70	0.56
23:DB:175:G:H2'	23:DB:176:A:C8	2.40	0.56
23:DB:2025:C:OP1	26:DD:154:LYS:HE3	2.04	0.56
23:DB:2061:G:H5''	23:DB:2503:A:C2	2.40	0.56
23:DB:241:A:O3'	23:DB:242:G:H4'	2.05	0.56
23:DB:467:G:O2'	23:DB:468:G:H5'	2.05	0.56
23:DB:479:A:H4'	23:DB:479:A:OP1	2.04	0.56
23:DB:552:U:O2'	23:DB:553:G:H5'	2.05	0.56
26:DD:29:VAL:HG22	26:DD:30:GLU:N	2.13	0.56
26:DD:69:ALA:HB1	26:DD:92:VAL:CG1	2.33	0.56
27:DE:149:ILE:HD13	27:DE:186:VAL:CG1	2.32	0.56
27:DE:17:THR:HG23	27:DE:18:THR:H	1.67	0.56
28:DF:163:GLU:HG2	28:DF:166:ARG:CZ	2.35	0.56
34:DM:6:ARG:HD2	34:DM:71:LYS:HE3	1.86	0.56
35:DN:103:ARG:HG2	35:DN:106:ASP:OD2	2.05	0.56
36:DO:108:ASP:O	36:DO:112:GLU:HG3	2.05	0.56
43:DW:45:HIS:HB2	43:DW:66:VAL:HG11	1.87	0.56
1:AA:1452:C:H5'	1:AA:1453:G:C4	2.40	0.56
1:AA:21:G:H2'	1:AA:22:G:C8	2.40	0.56
1:AA:321:A:O2'	1:AA:322:C:H5'	2.05	0.56
4:AE:104:ILE:HD11	4:AE:114:LEU:HB2	1.87	0.56
4:AE:114:LEU:HD13	4:AE:122:VAL:HG21	1.88	0.56
4:AE:96:GLN:HB3	4:AE:123:LEU:CD1	2.36	0.56
4:AE:45:VAL:O	4:AE:71:ILE:HG22	2.04	0.56
12:AM:53:ASP:OD1	12:AM:56:ARG:HD2	2.05	0.56
50:B3:4:LYS:NZ	50:B3:59:ALA:H	2.03	0.56
23:BB:1381:G:O2'	23:BB:1382:G:H5'	2.04	0.56
23:BB:2297:A:N6	23:BB:2319:G:C3'	2.59	0.56
23:BB:2590:A:H2'	23:BB:2591:C:C6	2.40	0.56
23:BB:2813:A:H2'	23:BB:2814:A:H8	1.70	0.56
27:BE:74:LYS:O	27:BE:82:GLY:HA3	2.05	0.56
27:BE:6:LYS:HG3	27:BE:9:GLN:HB2	1.86	0.56
23:BB:1081:U:C5'	52:BI:126:ARG:HD2	2.35	0.56
31:BJ:128:ASN:CG	31:BJ:129:GLU:H	2.08	0.56
32:BK:5:GLN:HA	32:BK:20:MET:SD	2.46	0.56
35:BN:46:ARG:O	35:BN:50:PRO:HG2	2.06	0.56
1:CA:1329:A:O2'	1:CA:1330:U:H5'	2.05	0.56
5:CF:2:ARG:HH11	5:CF:2:ARG:HB3	1.71	0.56
6:CG:29:LEU:HD23	6:CG:104:VAL:HG13	1.87	0.56
9:CJ:30:LYS:O	9:CJ:30:LYS:HD3	2.04	0.56
9:CJ:9:ARG:HB2	9:CJ:99:GLN:H	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CM:15:VAL:O	12:CM:19:THR:HG23	2.05	0.56
13:CN:29:ILE:HG22	13:CN:30:ILE:N	2.20	0.56
2:CC:39:ARG:NH2	13:CN:91:GLU:HB3	2.19	0.56
23:DB:2285:C:H5	48:D1:7:LYS:HE3	1.70	0.56
23:DB:1061:U:O4'	23:DB:1070:A:H1'	2.04	0.56
23:DB:2150:C:H2'	23:DB:2151:U:C6	2.39	0.56
23:DB:2250:G:N2	23:DB:2496:C:H4'	2.20	0.56
23:DB:2813:A:H2'	23:DB:2814:A:H8	1.70	0.56
23:DB:609:A:H2'	23:DB:610:C:O4'	2.05	0.56
23:DB:660:C:H2'	23:DB:661:A:C8	2.39	0.56
23:DB:69:C:H2'	23:DB:70:G:C8	2.40	0.56
25:DC:208:GLY:HA2	25:DC:212:TRP:CB	2.35	0.56
27:DE:126:VAL:HG21	27:DE:132:LYS:NZ	2.19	0.56
28:DF:4:HIS:O	28:DF:7:TYR:HB3	2.05	0.56
29:DG:6:ALA:HB3	29:DG:7:PRO:HD3	1.87	0.56
30:DH:37:VAL:O	30:DH:38:PRO:C	2.43	0.56
26:DD:156:PHE:CD1	31:DJ:81:ILE:HG21	2.41	0.56
39:DR:76:LYS:HA	39:DR:91:GLN:H	1.70	0.56
41:DT:76:ARG:HG2	41:DT:77:ARG:N	2.20	0.56
22:DA:75:G:H5''	24:DV:12:GLN:OE1	2.05	0.56
1:AA:218:U:H2'	1:AA:219:U:C6	2.40	0.56
1:AA:429:U:H1'	1:AA:430:A:H5''	1.87	0.56
1:AA:477:C:H2'	1:AA:478:A:C8	2.41	0.56
5:AF:55:HIS:O	5:AF:56:LYS:HG3	2.06	0.56
9:AJ:55:PRO:O	9:AJ:56:HIS:HB3	2.06	0.56
9:AJ:53:ILE:HG23	13:AN:84:ARG:CZ	2.36	0.56
33:BL:63:LYS:N	50:B3:24:LYS:HZ2	2.02	0.56
50:B3:44:ARG:CG	50:B3:45:PRO:HD2	2.34	0.56
23:BB:1076:C:H4'	52:BI:94:LYS:HZ2	1.70	0.56
23:BB:1082:U:C2	23:BB:1086:A:C6	2.93	0.56
23:BB:1250:G:H5''	38:BQ:5:ARG:HG2	1.86	0.56
23:BB:125:A:H4'	49:B2:13:ASN:OD1	2.05	0.56
23:BB:1528:A:H2'	23:BB:1529:G:O4'	2.06	0.56
23:BB:1695:G:H1'	25:BC:14:HIS:O	2.05	0.56
23:BB:796:C:H2'	23:BB:797:G:C8	2.40	0.56
27:BE:120:VAL:N	27:BE:189:THR:HG22	2.20	0.56
23:BB:328:U:H4'	42:BU:65:GLN:CG	2.35	0.56
24:BV:9:ARG:HH12	24:BV:27:PRO:HA	1.69	0.56
1:CA:266:G:O2'	1:CA:267:C:H3'	2.06	0.56
1:CA:586:C:C2'	1:CA:587:G:H5'	2.36	0.56
1:CA:981:U:H2'	1:CA:982:U:C5	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:162:GLU:HG3	3:CD:163:GLN:HE21	1.70	0.56
19:CT:57:VAL:HG12	19:CT:71:ALA:HB1	1.86	0.56
23:DB:1327:A:H2'	23:DB:1328:A:O4'	2.06	0.56
23:DB:131:A:H2'	23:DB:132:G:H8	1.70	0.56
23:DB:1709:U:H2'	23:DB:1710:G:C8	2.40	0.56
23:DB:17:G:H2'	23:DB:18:U:C6	2.40	0.56
23:DB:1826:G:H2'	23:DB:1827:U:H6	1.69	0.56
23:DB:2233:U:H2'	23:DB:2234:G:H8	1.69	0.56
23:DB:2872:A:O2'	23:DB:2873:A:H5''	2.05	0.56
23:DB:2882:A:H2'	23:DB:2883:A:H5''	1.86	0.56
23:DB:2898:U:H2'	23:DB:2899:A:H8	1.70	0.56
23:DB:351:C:H2'	23:DB:352:A:H8	1.69	0.56
23:DB:3:U:H2'	23:DB:4:U:H6	1.70	0.56
23:DB:490:C:H3'	23:DB:491:G:H5''	1.87	0.56
23:DB:527:C:O2	23:DB:527:C:O4'	2.21	0.56
26:DD:201:LEU:O	26:DD:202:ILE:HD12	2.05	0.56
27:DE:190:ALA:HB3	27:DE:193:VAL:HG22	1.86	0.56
27:DE:6:LYS:CD	27:DE:7:ASP:H	2.16	0.56
28:DF:163:GLU:HA	28:DF:166:ARG:HD2	1.87	0.56
28:DF:3:LEU:O	28:DF:3:LEU:HD23	2.05	0.56
28:DF:79:ARG:HB2	28:DF:82:TYR:CE2	2.40	0.56
37:DP:90:ALA:N	37:DP:112:ARG:HH21	2.01	0.56
1:AA:1040:U:H2'	1:AA:1041:G:C8	2.40	0.56
1:AA:384:G:H2'	1:AA:385:C:H6	1.69	0.56
1:AA:674:G:H2'	1:AA:675:A:C8	2.40	0.56
1:AA:677:U:H2'	1:AA:678:U:H6	1.70	0.56
1:AA:930:C:H2'	1:AA:931:C:O4'	2.05	0.56
2:AC:146:LYS:HE3	2:AC:202:PHE:CE2	2.38	0.56
5:AF:100:SER:HA	17:AR:23:LYS:NZ	2.19	0.56
1:AA:1506:U:H4'	10:AK:128:VAL:C	2.26	0.56
10:AK:25:SER:HB3	10:AK:28:ASN:O	2.06	0.56
12:AM:13:HIS:HB2	12:AM:16:ILE:CG2	2.36	0.56
13:AN:60:ARG:NH2	13:AN:62:ARG:HE	2.02	0.56
2:AC:33:ASP:HB2	13:AN:64:ARG:HD3	1.86	0.56
17:AR:41:SER:HB2	17:AR:51:GLN:HG2	1.87	0.56
21:AU:13:VAL:HG13	21:AU:14:ALA:N	2.12	0.56
48:B1:9:LYS:HG2	50:B3:34:LYS:HZ3	1.70	0.56
23:BB:1709:U:H2'	23:BB:1710:G:C8	2.41	0.56
23:BB:1729:U:H2'	23:BB:1730:C:H4'	1.87	0.56
23:BB:2869:G:H2'	23:BB:2870:C:C6	2.40	0.56
25:BC:174:ARG:HG3	25:BC:180:MET:SD	2.44	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:68:ARG:NH2	25:BC:190:THR:HG23	2.20	0.56
26:BD:201:LEU:N	26:BD:201:LEU:HD12	2.19	0.56
26:BD:5:VAL:HB	26:BD:27:ILE:O	2.05	0.56
27:BE:148:ILE:H	27:BE:183:PHE:HB3	1.70	0.56
27:BE:28:VAL:O	27:BE:32:VAL:HG23	2.06	0.56
28:BF:116:LEU:HD12	28:BF:176:PHE:CD2	2.40	0.56
28:BF:170:ALA:HA	28:BF:173:ASP:OD1	2.04	0.56
29:BG:132:LEU:HD11	29:BG:144:ALA:HB2	1.88	0.56
31:BJ:18:VAL:CG2	31:BJ:56:VAL:HA	2.35	0.56
32:BK:71:ARG:O	32:BK:72:PRO:C	2.44	0.56
34:BM:3:GLN:N	34:BM:4:PRO:CD	2.63	0.56
38:BQ:65:ASN:O	38:BQ:69:ARG:HG2	2.06	0.56
41:BT:68:LYS:NZ	41:BT:73:ARG:HA	2.20	0.56
1:CA:477:C:H2'	1:CA:478:A:C8	2.41	0.56
3:CD:192:ALA:HB3	3:CD:194:ILE:HG22	1.87	0.56
6:CG:55:LYS:HG3	6:CG:56:SER:H	1.70	0.56
8:CI:78:ILE:O	8:CI:82:ILE:HG13	2.04	0.56
12:CM:106:ARG:N	12:CM:106:ARG:HH11	2.04	0.56
23:DB:1381:G:H2'	23:DB:1382:G:H5'	1.86	0.56
23:DB:155:A:H2'	23:DB:156:A:C8	2.41	0.56
23:DB:1661:G:O2'	23:DB:1662:U:H5'	2.05	0.56
23:DB:171:U:H2'	23:DB:172:A:C8	2.40	0.56
23:DB:172:A:H2'	23:DB:173:A:H8	1.70	0.56
23:DB:1771:C:H2'	23:DB:1772:A:H8	1.69	0.56
23:DB:1917:U:O2'	23:DB:1918:A:H5'	2.06	0.56
23:DB:2140:G:H2'	23:DB:2141:G:O4'	2.04	0.56
23:DB:234:U:O2'	23:DB:235:U:H5'	2.04	0.56
23:DB:2626:C:O2'	23:DB:2627:G:H5'	2.05	0.56
23:DB:877:A:N6	23:DB:898:C:H2'	2.19	0.56
23:DB:91:A:H1'	23:DB:92:U:C6	2.41	0.56
25:DC:79:ARG:HD2	25:DC:110:LYS:HE2	1.87	0.56
31:DJ:25:LEU:N	31:DJ:25:LEU:HD13	2.20	0.56
32:DK:113:MET:CE	32:DK:116:ILE:HD11	2.35	0.56
33:DL:92:LEU:HD13	33:DL:96:LYS:HG2	1.87	0.56
35:DN:86:ARG:NH2	35:DN:117:ASP:HA	2.20	0.56
35:DN:4:ARG:HG2	35:DN:5:LYS:N	2.19	0.56
26:DD:8:LYS:CE	37:DP:5:LYS:HG3	2.34	0.56
39:DR:67:GLY:H	39:DR:98:ILE:HA	1.69	0.56
23:DB:975:A:H5'	39:DR:83:TYR:OH	2.05	0.56
40:DS:28:LYS:HA	40:DS:31:GLN:NE2	2.19	0.56
40:DS:46:LEU:O	40:DS:50:VAL:HG13	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:301:G:H2'	1:AA:302:G:H8	1.71	0.56
1:AA:499:A:H4'	1:AA:500:G:OP1	2.06	0.56
1:AA:77:A:H2'	1:AA:78:A:C8	2.40	0.56
20:AB:195:VAL:HG12	20:AB:197:PHE:H	1.70	0.56
20:AB:76:SER:HA	20:AB:92:ASN:HB2	1.87	0.56
6:AG:114:SER:C	6:AG:118:ARG:HG3	2.26	0.56
19:AT:82:ILE:HG13	19:AT:83:ASN:N	2.20	0.56
21:AU:38:GLU:C	21:AU:40:PRO:HD2	2.26	0.56
48:B1:53:ILE:HD13	48:B1:54:LYS:H	1.70	0.56
23:BB:682:G:H5'	49:B2:26:ASN:CG	2.25	0.56
23:BB:458:G:H5''	49:B2:39:ARG:O	2.04	0.56
50:B3:34:LYS:C	50:B3:35:LYS:HD3	2.25	0.56
50:B3:41:ARG:NH1	50:B3:42:HIS:HB3	2.20	0.56
22:BA:26:C:H2'	22:BA:27:C:O4'	2.06	0.56
23:BB:1061:U:O4'	23:BB:1070:A:H1'	2.05	0.56
23:BB:1351:C:O2'	23:BB:1571:A:H1'	2.06	0.56
23:BB:1459:G:C2'	23:BB:1460:U:H5'	2.35	0.56
23:BB:189:G:OP1	46:BZ:24:ILE:HD12	2.05	0.56
23:BB:1938:A:O2'	23:BB:1939:U:H5''	2.06	0.56
23:BB:2022:U:O2'	23:BB:2617:U:H5'	2.06	0.56
23:BB:2147:A:H5'	23:BB:2148:G:H4'	1.87	0.56
23:BB:962:G:H21	23:BB:2250:G:H1	1.53	0.56
23:BB:2053:G:H5''	26:BD:150:GLN:HA	1.86	0.56
28:BF:83:PRO:HG2	28:BF:84:ILE:H	1.71	0.56
30:BH:80:ILE:HD12	30:BH:102:ALA:HB2	1.88	0.56
31:BJ:58:ASN:HA	31:BJ:127:GLY:N	2.21	0.56
31:BJ:16:TYR:O	31:BJ:55:ILE:HG22	2.04	0.56
31:BJ:7:LYS:HZ3	31:BJ:48:VAL:HB	1.70	0.56
33:BL:80:SER:O	33:BL:116:VAL:HA	2.06	0.56
33:BL:119:PRO:O	33:BL:119:PRO:HD2	2.05	0.56
34:BM:33:LEU:HB2	34:BM:101:VAL:CG2	2.35	0.56
38:BQ:34:ALA:O	38:BQ:38:VAL:HG12	2.06	0.56
38:BQ:79:ILE:HA	38:BQ:82:LEU:HD12	1.87	0.56
39:BR:85:LYS:C	39:BR:87:GLN:H	2.09	0.56
39:BR:66:HIS:H	39:BR:98:ILE:HD13	1.70	0.56
1:CA:1314:C:OP2	18:CS:5:LYS:HG2	2.05	0.56
1:CA:1470:U:O2'	1:CA:1471:U:H5'	2.05	0.56
1:CA:474:G:H2'	1:CA:475:C:C6	2.40	0.56
1:CA:815:A:H62	1:CA:1509:C:H1'	1.70	0.56
20:CB:78:ALA:HB1	20:CB:213:LEU:HD13	1.87	0.56
6:CG:49:LEU:C	6:CG:51:GLN:H	2.07	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CK:22:ILE:HD13	10:CK:95:THR:HG23	1.87	0.56
15:CP:52:LEU:HD21	15:CP:74:LEU:HB2	1.87	0.56
50:D3:4:LYS:HE3	50:D3:61:LEU:HD13	1.88	0.56
23:DB:144:A:H8	23:DB:144:A:O5'	1.89	0.56
23:DB:2090:A:O2'	46:DZ:49:ARG:CD	2.53	0.56
23:DB:876:C:H3'	23:DB:877:A:O4'	2.06	0.56
25:DC:128:THR:HA	25:DC:190:THR:HA	1.86	0.56
25:DC:179:GLU:HB3	25:DC:266:ILE:HA	1.88	0.56
26:DD:56:LYS:HB3	26:DD:56:LYS:HZ2	1.69	0.56
23:DB:323:C:C1'	27:DE:164:LEU:HB3	2.25	0.56
28:DF:174:PHE:H	28:DF:175:PRO:CD	2.19	0.56
52:DI:24:GLY:HA2	52:DI:34:ILE:HD12	1.86	0.56
42:DU:25:LYS:O	42:DU:34:ILE:HG12	2.06	0.56
23:DB:2262:U:OP2	43:DW:13:ARG:HD3	2.04	0.56
43:DW:38:ARG:HG3	43:DW:39:GLN:N	2.21	0.56
1:AA:1225:A:H3'	1:AA:1226:C:C6	2.40	0.56
1:AA:1329:A:O2'	1:AA:1330:U:H5'	2.05	0.56
1:AA:1463:U:H2'	1:AA:1464:U:C6	2.40	0.56
1:AA:1470:U:O2'	1:AA:1471:U:H5'	2.06	0.56
1:AA:474:G:H2'	1:AA:475:C:C6	2.40	0.56
1:AA:586:C:C2'	1:AA:587:G:H5'	2.35	0.56
20:AB:163:ILE:CG2	20:AB:164:ASP:H	1.96	0.56
3:AD:117:VAL:O	3:AD:130:ASN:HA	2.05	0.56
6:AG:22:LEU:O	6:AG:25:PHE:HB3	2.06	0.56
6:AG:91:ARG:HB3	6:AG:92:PRO:HD2	1.88	0.56
14:AO:63:ARG:HA	14:AO:66:LEU:HD12	1.87	0.56
18:AS:48:ILE:N	18:AS:48:ILE:HD12	2.21	0.56
19:AT:61:ALA:HA	19:AT:66:ILE:O	2.05	0.56
23:BB:1661:G:O2'	23:BB:1662:U:H5'	2.05	0.56
23:BB:1826:G:H2'	23:BB:1827:U:H6	1.70	0.56
23:BB:1842:G:H2'	23:BB:1843:C:H6	1.70	0.56
23:BB:1889:A:H2'	23:BB:1890:A:C8	2.41	0.56
23:BB:2729:G:H2'	23:BB:2730:C:C6	2.41	0.56
23:BB:2747:G:O6	23:BB:2755:C:H5''	2.05	0.56
23:BB:2840:C:H5''	35:BN:53:THR:CG2	2.36	0.56
23:BB:2:G:O2'	23:BB:3:U:H5'	2.05	0.56
23:BB:597:G:H4'	33:BL:21:ARG:NE	2.18	0.56
23:BB:851:C:O2'	45:BY:45:GLY:HA3	2.05	0.56
29:BG:70:LEU:HD12	29:BG:74:MET:HE3	1.87	0.56
52:BI:10:LEU:HD12	52:BI:10:LEU:O	2.05	0.56
33:BL:118:THR:N	33:BL:119:PRO:CD	2.66	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BM:41:LEU:HG	34:BM:95:LEU:HD12	1.86	0.56
34:BM:96:ILE:HD13	34:BM:97:GLN:N	2.21	0.56
36:BO:8:ILE:HG23	36:BO:9:ARG:N	2.20	0.56
37:BP:47:ILE:HG23	37:BP:63:ILE:HG13	1.88	0.56
37:BP:80:VAL:O	37:BP:80:VAL:HG13	2.05	0.56
38:BQ:22:GLY:O	38:BQ:24:TYR:N	2.39	0.56
38:BQ:50:ARG:O	38:BQ:53:LYS:HD3	2.05	0.56
23:BB:535:G:O2'	38:BQ:52:ARG:HD3	2.05	0.56
41:BT:96:VAL:HG22	41:BT:97:GLY:N	2.21	0.56
1:CA:218:U:H2'	1:CA:219:U:C6	2.40	0.56
1:CA:429:U:H1'	1:CA:430:A:H5''	1.87	0.56
1:CA:764:C:H2'	1:CA:765:G:H5'	1.88	0.56
6:CG:2:ARG:HH11	6:CG:2:ARG:CB	2.17	0.56
9:CJ:9:ARG:HB3	9:CJ:99:GLN:HE21	1.70	0.56
1:CA:1228:C:H5'	12:CM:112:ARG:HB3	1.87	0.56
1:CA:1313:U:P	18:CS:5:LYS:HA	2.46	0.56
23:DB:2527:C:H1'	51:D4:1:MET:HG3	1.87	0.56
23:DB:1316:U:O2'	23:DB:1317:G:H5'	2.05	0.56
23:DB:2600:A:O2'	23:DB:2601:C:H5'	2.05	0.56
23:DB:2729:G:H2'	23:DB:2730:C:C6	2.41	0.56
23:DB:2758:A:O2'	23:DB:2759:G:H5'	2.04	0.56
23:DB:2798:U:H1'	23:DB:2800:A:N6	2.20	0.56
23:DB:287:G:H2'	23:DB:288:U:C6	2.40	0.56
23:DB:818:G:H5'	23:DB:839:U:OP1	2.06	0.56
25:DC:156:SER:O	25:DC:158:GLY:N	2.39	0.56
26:DD:159:LYS:O	26:DD:160:LYS:HB2	2.05	0.56
31:DJ:7:LYS:HD2	31:DJ:45:THR:CB	2.35	0.56
27:DE:184:ASP:CB	33:DL:13:LYS:HE3	2.36	0.56
23:DB:2360:G:O2'	33:DL:61:LEU:HD21	2.06	0.56
36:DO:15:ARG:NH1	43:DW:76:ARG:HE	2.03	0.56
37:DP:28:LYS:HD3	37:DP:44:GLY:CA	2.33	0.56
41:DT:34:VAL:HG21	41:DT:43:ILE:CD1	2.33	0.56
42:DU:21:ARG:HG3	42:DU:21:ARG:HH11	1.70	0.56
46:DZ:41:HIS:ND1	46:DZ:42:PRO:HD2	2.20	0.56
1:AA:1310:G:H2'	1:AA:1311:A:O4'	2.06	0.56
1:AA:1468:A:O2'	1:AA:1469:C:H5'	2.05	0.56
1:AA:327:A:H1'	1:AA:329:A:O4'	2.06	0.56
1:AA:956:U:O2'	1:AA:957:U:H5'	2.06	0.56
20:AB:144:GLU:O	20:AB:148:GLY:HA3	2.05	0.56
20:AB:53:LEU:HD21	20:AB:212:TYR:HE2	1.70	0.56
3:AD:10:LEU:HD21	3:AD:62:ARG:CD	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:38:PHE:O	8:AI:41:GLU:HB2	2.06	0.56
8:AI:99:LYS:C	8:AI:99:LYS:HD3	2.26	0.56
10:AK:92:ARG:CB	10:AK:92:ARG:HH11	2.19	0.56
22:BA:90:C:H2'	22:BA:91:C:O4'	2.05	0.56
23:BB:1133:A:H4'	23:BB:1134:A:H5''	1.87	0.56
23:BB:988:A:H4'	23:BB:1155:A:N1	2.21	0.56
23:BB:1750:G:O2'	23:BB:1751:U:H5'	2.06	0.56
23:BB:2239:G:H5'	25:BC:246:PRO:HD2	1.87	0.56
23:BB:2586:U:H2'	23:BB:2587:A:C8	2.41	0.56
23:BB:2665:A:O2'	23:BB:2666:C:H5'	2.06	0.56
23:BB:2798:U:H1'	23:BB:2800:A:N6	2.20	0.56
23:BB:2052:A:H5'	26:BD:148:GLN:O	2.05	0.56
27:BE:128:ALA:O	27:BE:157:LEU:HB2	2.05	0.56
29:BG:159:LYS:HG3	29:BG:159:LYS:O	2.05	0.56
31:BJ:102:GLU:HG3	31:BJ:124:VAL:HG11	1.87	0.56
1:AA:1432:G:H5''	37:BP:106:ALA:HB2	1.88	0.56
23:BB:1754:A:P	37:BP:94:ALA:HB1	2.46	0.56
39:BR:90:ARG:HH22	39:BR:92:TRP:HB3	1.69	0.56
39:BR:66:HIS:HA	39:BR:98:ILE:HD13	1.86	0.56
40:BS:9:HIS:O	40:BS:100:THR:HB	2.04	0.56
40:BS:47:VAL:O	40:BS:51:LEU:HB2	2.05	0.56
40:BS:8:ARG:CZ	40:BS:102:HIS:NE2	2.69	0.56
41:BT:40:LYS:HB3	41:BT:58:VAL:HB	1.88	0.56
41:BT:30:ILE:HG22	41:BT:84:TYR:O	2.06	0.56
45:BY:13:ILE:HG23	45:BY:19:HIS:HE1	1.70	0.56
1:CA:1262:C:H2'	1:CA:1263:C:H6	1.70	0.56
3:CD:106:PHE:CD1	3:CD:158:LEU:HD21	2.41	0.56
1:CA:642:A:C5	7:CH:106:SER:HA	2.41	0.56
1:CA:1250:A:H5''	8:CI:69:GLY:N	2.21	0.56
10:CK:78:ILE:CD1	10:CK:78:ILE:H	2.06	0.56
23:DB:1018:U:O2'	23:DB:1019:U:H5'	2.05	0.56
23:DB:1528:A:H2'	23:DB:1529:G:O4'	2.06	0.56
23:DB:171:U:H2'	23:DB:172:A:H8	1.71	0.56
23:DB:1750:G:O2'	23:DB:1751:U:H5'	2.06	0.56
23:DB:388:G:N7	23:DB:390:U:H2'	2.20	0.56
25:DC:136:VAL:C	25:DC:138:SER:H	2.09	0.56
23:DB:1796:U:OP1	25:DC:269:ARG:HD3	2.06	0.56
27:DE:17:THR:HG23	27:DE:18:THR:N	2.21	0.56
27:DE:195:GLN:HB2	27:DE:198:GLU:O	2.06	0.56
23:DB:659:G:H21	27:DE:30:GLN:NE2	2.03	0.56
27:DE:85:PHE:O	27:DE:86:ALA:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:DH:58:LEU:O	30:DH:61:VAL:HG22	2.05	0.56
30:DH:82:SER:O	30:DH:83:LYS:HD2	2.05	0.56
33:DL:78:ARG:CZ	33:DL:113:ALA:HB1	2.36	0.56
33:DL:132:ARG:NH1	33:DL:140:GLY:HA2	2.21	0.56
24:DV:9:ARG:NH2	24:DV:20:LEU:HD11	2.20	0.56
45:DY:4:ILE:HG12	45:DY:5:LYS:HG3	1.87	0.56
1:AA:449:G:H2'	1:AA:450:G:C8	2.40	0.56
1:AA:552:U:H2'	1:AA:553:A:C8	2.41	0.56
4:AE:87:VAL:HG11	4:AE:92:ARG:HD3	1.87	0.56
11:AL:106:VAL:HG21	11:AL:109:ARG:HG3	1.88	0.56
23:BB:1996:C:H5	32:BK:32:TYR:HH	1.53	0.56
23:BB:2065:C:H2'	23:BB:2066:C:H6	1.71	0.56
23:BB:2262:U:H2'	23:BB:2263:C:H6	1.71	0.56
23:BB:2484:G:H1'	34:BM:119:LEU:HD12	1.87	0.56
23:BB:1971:U:H1'	25:BC:238:ASN:ND2	2.20	0.56
26:BD:8:LYS:HA	26:BD:197:THR:HG23	1.88	0.56
28:BF:7:TYR:O	28:BF:11:VAL:HB	2.05	0.56
28:BF:66:ILE:HA	28:BF:85:GLY:O	2.06	0.56
29:BG:84:LYS:HA	29:BG:84:LYS:HZ3	1.70	0.56
30:BH:69:ALA:HB2	30:BH:140:ALA:HA	1.86	0.56
34:BM:8:LYS:HB2	34:BM:12:MET:HE3	1.87	0.56
35:BN:32:GLU:HB3	35:BN:33:ILE:HD13	1.88	0.56
35:BN:73:ASN:N	35:BN:73:ASN:HD22	2.03	0.56
39:BR:42:ALA:HB1	39:BR:45:GLU:OE2	2.05	0.56
41:BT:30:ILE:HG23	41:BT:32:LEU:HD13	1.87	0.56
43:BW:39:GLN:HG2	43:BW:68:PHE:HB2	1.86	0.56
1:CA:321:A:O2'	1:CA:322:C:H5'	2.06	0.56
1:CA:327:A:H1'	1:CA:329:A:O4'	2.06	0.56
1:CA:591:U:H2'	1:CA:592:G:H8	1.70	0.56
5:CF:47:LEU:HD22	17:CR:65:SER:OG	2.05	0.56
6:CG:24:LYS:O	6:CG:28:ILE:HG12	2.06	0.56
11:CL:107:LYS:O	11:CL:108:ASP:HB2	2.06	0.56
13:CN:41:TRP:O	13:CN:45:LEU:HG	2.06	0.56
23:DB:1862:G:O2'	23:DB:1863:G:H5'	2.06	0.56
23:DB:1878:G:H2'	23:DB:1879:C:H6	1.69	0.56
23:DB:2639:A:H2'	23:DB:2640:G:O4'	2.06	0.56
23:DB:2839:G:H2'	23:DB:2840:C:H6	1.69	0.56
23:DB:374:A:H61	23:DB:400:G:H1'	1.68	0.56
23:DB:4:U:O2'	23:DB:5:A:H5'	2.05	0.56
25:DC:224:MET:HA	25:DC:233:GLY:N	2.13	0.56
25:DC:254:LYS:HD2	25:DC:254:LYS:N	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:25:THR:HG22	26:DD:188:LEU:HG	1.88	0.56
26:DD:35:THR:HB	26:DD:48:ILE:HG13	1.87	0.56
28:DF:103:ILE:HD12	28:DF:104:THR:N	2.21	0.56
29:DG:125:PRO:HG2	29:DG:129:GLU:HB2	1.88	0.56
32:DK:69:VAL:HG11	32:DK:106:GLU:CD	2.26	0.56
23:DB:2405:G:H5''	33:DL:70:LYS:HG3	1.86	0.56
39:DR:6:GLN:HB3	39:DR:41:ILE:CD1	2.36	0.56
43:DW:35:ILE:HB	43:DW:67:LYS:NZ	2.20	0.56
1:AA:1329:A:OP1	12:AM:28:ARG:HB2	2.05	0.56
1:AA:529:G:H22	11:AL:47:ALA:CB	2.18	0.56
13:AN:60:ARG:HE	13:AN:62:ARG:HE	1.53	0.56
15:AP:20:VAL:HG22	15:AP:21:VAL:N	2.21	0.56
1:AA:332:G:OP2	19:AT:2:ASN:HB3	2.06	0.56
48:B1:42:VAL:HG12	48:B1:43:ARG:HE	1.71	0.56
23:BB:100:U:O3'	42:BU:90:LYS:HG2	2.06	0.56
23:BB:1739:A:H2'	23:BB:1740:G:C8	2.40	0.56
23:BB:1923:U:H2'	23:BB:1924:C:C6	2.41	0.56
23:BB:758:C:O2	23:BB:1981:A:H2	1.86	0.56
23:BB:198:C:C2'	23:BB:199:A:H5''	2.35	0.56
23:BB:2239:G:H5'	25:BC:246:PRO:CG	2.35	0.56
23:BB:2771:C:H2'	23:BB:2772:C:C6	2.41	0.56
23:BB:765:C:H2'	23:BB:766:U:C6	2.39	0.56
23:BB:974:G:O4'	23:BB:990:A:N6	2.39	0.56
27:BE:149:ILE:CG1	27:BE:186:VAL:HG13	2.36	0.56
29:BG:88:LEU:HD11	29:BG:95:ALA:HB2	1.88	0.56
30:BH:140:ALA:C	30:BH:141:LYS:HD3	2.26	0.56
30:BH:128:HIS:CB	30:BH:144:VAL:HB	2.35	0.56
32:BK:10:VAL:HG11	32:BK:15:GLY:O	2.04	0.56
33:BL:110:VAL:HB	33:BL:126:ARG:CZ	2.35	0.56
38:BQ:35:PHE:HE1	39:BR:85:LYS:HZ2	1.53	0.56
40:BS:27:LYS:CD	40:BS:27:LYS:H	2.11	0.56
40:BS:89:ALA:C	40:BS:90:LYS:HZ2	2.08	0.56
44:BX:45:GLN:N	44:BX:45:GLN:NE2	2.53	0.56
1:CA:1292:G:H2'	1:CA:1293:C:H6	1.71	0.56
1:CA:235:C:H2'	1:CA:236:A:H8	1.70	0.56
1:CA:552:U:H2'	1:CA:553:A:C8	2.41	0.56
1:CA:93:U:C2'	1:CA:94:G:H5'	2.36	0.56
2:CC:35:ASP:O	2:CC:39:ARG:HG3	2.06	0.56
3:CD:22:SER:H	3:CD:109:THR:CG2	2.19	0.56
4:CE:55:VAL:N	4:CE:56:PRO:HD2	2.21	0.56
4:CE:67:ARG:HB2	4:CE:67:ARG:NH1	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CH:100:ILE:HG13	7:CH:128:VAL:CG2	2.36	0.56
7:CH:35:ILE:HG22	7:CH:39:LEU:HD23	1.88	0.56
13:CN:12:ARG:HG3	13:CN:53:ASP:O	2.05	0.56
1:CA:617:G:H5'	15:CP:46:LYS:HD3	1.87	0.56
16:CQ:10:ARG:NH1	16:CQ:55:GLY:H	2.04	0.56
18:CS:11:ASP:O	18:CS:14:LEU:HG	2.06	0.56
47:D0:47:TYR:HB2	47:D0:50:GLY:O	2.06	0.56
23:DB:1082:U:N3	23:DB:1086:A:C2	2.73	0.56
23:DB:1438:U:N3	23:DB:1552:A:N6	2.53	0.56
23:DB:1854:A:N6	23:DB:1888:G:H1'	2.21	0.56
23:DB:1889:A:H2'	23:DB:1890:A:C8	2.41	0.56
23:DB:2443:C:O2'	23:DB:2444:G:H5'	2.05	0.56
23:DB:2771:C:H2'	23:DB:2772:C:C6	2.41	0.56
23:DB:458:G:H22	23:DB:469:G:H2'	1.68	0.56
23:DB:581:C:OP1	38:DQ:31:TYR:HB3	2.06	0.56
23:DB:871:U:H2'	23:DB:872:U:H6	1.70	0.56
25:DC:145:MET:HG2	25:DC:152:GLN:HG2	1.88	0.56
25:DC:68:ARG:HD3	25:DC:103:ILE:HD12	1.88	0.56
26:DD:33:ARG:HB2	26:DD:33:ARG:NH1	2.18	0.56
26:DD:50:VAL:HG11	26:DD:75:ALA:HB3	1.87	0.56
27:DE:163:ASN:HB2	27:DE:167:VAL:O	2.05	0.56
27:DE:191:ASP:O	27:DE:194:LYS:HG2	2.06	0.56
29:DG:94:ARG:HB2	29:DG:105:SER:CB	2.35	0.56
23:DB:1098:A:C3'	52:DI:3:LYS:HA	2.24	0.56
31:DJ:100:VAL:HG22	31:DJ:101:ILE:N	2.21	0.56
33:DL:76:GLU:OE1	33:DL:76:GLU:HA	2.05	0.56
37:DP:59:THR:HG22	37:DP:74:GLN:HB3	1.88	0.56
38:DQ:47:ARG:O	38:DQ:51:GLN:HG3	2.06	0.56
39:DR:40:MET:O	39:DR:54:VAL:HG22	2.06	0.56
1:AA:1090:U:H2'	1:AA:1091:U:C6	2.40	0.56
1:AA:69:G:H2'	1:AA:70:U:C6	2.40	0.56
1:AA:824:G:H2'	1:AA:825:A:H8	1.71	0.56
1:AA:97:G:H2'	1:AA:98:A:O4'	2.05	0.56
2:AC:26:LYS:HD3	2:AC:26:LYS:C	2.25	0.56
6:AG:45:ALA:HB2	6:AG:116:ALA:O	2.05	0.56
7:AH:111:THR:HG23	7:AH:114:ALA:HB2	1.87	0.56
8:AI:18:VAL:HG21	8:AI:82:ILE:HG13	1.88	0.56
8:AI:64:ILE:HD13	8:AI:78:ILE:HG21	1.87	0.56
10:AK:12:ARG:N	10:AK:12:ARG:HD2	2.21	0.56
17:AR:49:LYS:HA	17:AR:52:ARG:HD2	1.88	0.56
48:B1:31:GLU:HB3	48:B1:32:LYS:HE3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1460:U:H5''	23:BB:1461:C:C5	2.41	0.56
23:BB:170:U:H2'	23:BB:171:U:H6	1.71	0.56
23:BB:1862:G:O2'	23:BB:1863:G:H5'	2.06	0.56
23:BB:2825:G:H2'	23:BB:2826:A:H5'	1.88	0.56
27:BE:122:GLU:HB3	27:BE:152:GLU:CB	2.35	0.56
31:BJ:35:ARG:HH21	31:BJ:40:HIS:N	2.04	0.56
31:BJ:10:THR:HG22	31:BJ:50:THR:CA	2.35	0.56
23:BB:943:A:OP1	33:BL:42:SER:HB3	2.06	0.56
34:BM:108:VAL:N	34:BM:109:PRO:CD	2.69	0.56
40:BS:7:HIS:CG	40:BS:8:ARG:H	2.23	0.56
42:BU:23:LYS:NZ	42:BU:24:VAL:HG22	2.21	0.56
43:BW:42:THR:CB	43:BW:66:VAL:H	2.19	0.56
46:BZ:15:SER:HB2	46:BZ:25:ARG:HH12	1.71	0.56
46:BZ:48:GLN:H	46:BZ:51:VAL:HB	1.69	0.56
1:CA:1002:G:H2'	1:CA:1003:G:O4'	2.06	0.56
1:CA:1063:C:H3'	1:CA:1064:G:H2'	1.87	0.56
1:CA:272:C:H2'	1:CA:273:U:H6	1.70	0.56
1:CA:923:A:OP1	4:CE:25:LYS:HB3	2.05	0.56
1:CA:948:C:O2'	1:CA:949:A:H5'	2.05	0.56
20:CB:31:PHE:HB2	20:CB:41:ASN:HA	1.88	0.56
3:CD:151:GLN:NE2	3:CD:153:ARG:HD2	2.20	0.56
1:CA:875:U:O2'	7:CH:14:ARG:HD2	2.05	0.56
12:CM:78:ARG:NH1	12:CM:79:LEU:HD23	2.19	0.56
13:CN:42:ASN:HB3	13:CN:46:LYS:HE2	1.86	0.56
18:CS:62:THR:HG21	18:CS:64:GLU:OE1	2.05	0.56
48:D1:24:LYS:HZ2	48:D1:25:ASN:N	2.04	0.56
23:DB:1459:G:C2'	23:DB:1460:U:H5'	2.35	0.56
23:DB:2562:U:H1'	32:DK:23:LYS:HE2	1.88	0.56
23:DB:2665:A:O2'	23:DB:2666:C:H5'	2.05	0.56
23:DB:2888:C:H2'	23:DB:2889:C:H6	1.69	0.56
23:DB:324:A:H2'	23:DB:325:G:O4'	2.06	0.56
23:DB:621:A:H2'	23:DB:622:G:O4'	2.06	0.56
23:DB:703:U:H2'	23:DB:704:G:O4'	2.05	0.56
26:DD:146:ILE:CD1	26:DD:155:VAL:HG13	2.36	0.56
23:DB:675:A:OP1	27:DE:58:LYS:HE2	2.06	0.56
28:DF:107:VAL:HG12	28:DF:108:PRO:CD	2.32	0.56
35:DN:72:ASP:OD1	35:DN:74:GLU:HB3	2.05	0.56
35:DN:87:PHE:HB2	35:DN:94:TYR:CE2	2.40	0.56
24:DV:6:ALA:O	24:DV:65:VAL:HA	2.06	0.56
43:DW:42:THR:N	43:DW:65:LYS:HA	2.20	0.56
44:DX:25:GLN:HG2	44:DX:26:PHE:N	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:129:A:H1'	1:AA:130:A:C8	2.41	0.56
1:AA:1422:G:OP1	32:BK:48:PRO:HA	2.06	0.56
1:AA:266:G:O2'	1:AA:267:C:H3'	2.06	0.56
4:AE:104:ILE:HG23	4:AE:111:ARG:NH1	2.21	0.56
6:AG:24:LYS:O	6:AG:28:ILE:HG13	2.06	0.56
6:AG:61:PHE:O	6:AG:65:LEU:HD13	2.06	0.56
7:AH:65:PHE:CD2	7:AH:66:GLN:HG3	2.41	0.56
18:AS:68:HIS:NE2	18:AS:72:GLU:HG3	2.21	0.56
47:B0:16:ARG:HA	47:B0:19:ASP:OD2	2.05	0.56
33:BL:63:LYS:HA	50:B3:24:LYS:HD3	1.87	0.56
50:B3:53:ASP:O	50:B3:56:LEU:HB3	2.06	0.56
23:BB:1064:C:H4'	52:BI:90:GLY:HA2	1.88	0.56
23:BB:1080:A:H2'	23:BB:1081:U:H6	1.71	0.56
23:BB:1316:U:O2'	23:BB:1317:G:H5'	2.05	0.56
23:BB:140:C:H4'	23:BB:141:G:H21	1.71	0.56
23:BB:1534:U:H2'	23:BB:1536:C:N3	2.21	0.56
23:BB:1550:C:H2'	23:BB:1551:A:C8	2.40	0.56
23:BB:1656:C:P	26:BD:141:ARG:HD3	2.45	0.56
23:BB:1710:G:H4'	23:BB:2858:C:O2	2.06	0.56
23:BB:2484:G:OP1	34:BM:44:ARG:HD2	2.06	0.56
23:BB:2626:C:O2'	23:BB:2627:G:H5'	2.06	0.56
23:BB:527:C:H5''	23:BB:2779:U:C2	2.41	0.56
23:BB:315:G:H2'	23:BB:316:C:C6	2.41	0.56
23:BB:581:C:H2'	23:BB:582:A:H8	1.70	0.56
23:BB:639:U:H2'	23:BB:640:C:H6	1.69	0.56
23:BB:996:A:O3'	38:BQ:92:LYS:HE2	2.06	0.56
26:BD:119:ALA:HA	26:BD:123:LYS:HZ1	1.69	0.56
32:BK:40:LYS:HE3	32:BK:57:VAL:HG12	1.88	0.56
23:BB:663:G:OP1	33:BL:27:LEU:HD21	2.06	0.56
34:BM:3:GLN:OE1	34:BM:47:GLU:HB2	2.05	0.56
34:BM:53:MET:O	34:BM:54:THR:HG23	2.06	0.56
37:BP:29:VAL:CG2	37:BP:47:ILE:HD11	2.36	0.56
1:CA:1234:C:H1'	1:CA:1364:U:O2	2.06	0.56
1:CA:426:U:H2'	1:CA:427:U:C6	2.41	0.56
1:CA:576:C:OP2	1:CA:576:C:H3'	2.06	0.56
1:CA:676:A:H5''	10:CK:114:PRO:HB2	1.87	0.56
11:CL:30:ARG:O	11:CL:56:LEU:HA	2.06	0.56
12:CM:3:ILE:HA	12:CM:56:ARG:NH1	2.21	0.56
16:CQ:24:ILE:N	16:CQ:24:ILE:HD12	2.21	0.56
1:CA:673:A:H1'	17:CR:63:TYR:CE1	2.40	0.56
18:CS:15:LEU:HA	18:CS:18:VAL:CG1	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:CT:61:ALA:HA	19:CT:67:HIS:N	2.17	0.56
48:D1:7:LYS:HB3	48:D1:24:LYS:NZ	2.21	0.56
48:D1:42:VAL:O	48:D1:43:ARG:HB2	2.06	0.56
48:D1:9:LYS:HA	48:D1:24:LYS:CB	2.36	0.56
50:D3:12:ARG:O	50:D3:13:PHE:HB2	2.06	0.56
23:DB:1360:G:C2'	23:DB:1361:G:H5'	2.36	0.56
23:DB:1534:U:H2'	23:DB:1536:C:N3	2.21	0.56
23:DB:2074:U:H2'	23:DB:2075:U:C6	2.41	0.56
23:DB:30:G:H2'	23:DB:31:C:H6	1.71	0.56
23:DB:566:U:O2'	23:DB:567:U:H5'	2.06	0.56
23:DB:581:C:H2'	23:DB:582:A:H8	1.70	0.56
25:DC:63:ILE:HG22	25:DC:64:VAL:N	2.21	0.56
23:DB:2636:C:C5'	26:DD:80:TRP:HE1	2.19	0.56
28:DF:107:VAL:N	28:DF:108:PRO:CD	2.69	0.56
28:DF:16:MET:O	28:DF:20:ASN:HA	2.06	0.56
23:DB:1666:G:O3'	32:DK:6:THR:HG23	2.06	0.56
33:DL:90:VAL:HG13	33:DL:90:VAL:O	2.06	0.56
34:DM:19:GLY:C	34:DM:20:LEU:HD12	2.25	0.56
39:DR:79:ARG:HD2	39:DR:80:ARG:N	2.21	0.56
42:DU:25:LYS:HA	42:DU:34:ILE:H	1.71	0.56
43:DW:21:GLY:HA3	43:DW:32:ALA:HA	1.86	0.56
46:DZ:11:GLU:H	46:DZ:27:THR:HG22	1.71	0.56
23:DB:2091:C:H5'	46:DZ:49:ARG:HG3	1.88	0.56
1:AA:1404:C:H2'	1:AA:1405:G:C8	2.41	0.55
1:AA:272:C:H2'	1:AA:273:U:H6	1.71	0.55
1:AA:383:A:H2'	1:AA:384:G:O4'	2.05	0.55
1:AA:958:A:N1	18:AS:53:GLY:HA3	2.21	0.55
20:AB:125:PHE:HD2	20:AB:125:PHE:H	1.53	0.55
2:AC:181:ILE:N	2:AC:181:ILE:HD12	2.20	0.55
3:AD:58:GLN:OE1	3:AD:62:ARG:HG3	2.07	0.55
6:AG:72:VAL:HA	6:AG:89:GLU:HA	1.86	0.55
1:AA:450:G:H4'	15:AP:41:PRO:HB2	1.87	0.55
16:AQ:4:ILE:CD1	16:AQ:5:ARG:H	2.20	0.55
16:AQ:8:GLN:HB3	16:AQ:57:VAL:HG13	1.88	0.55
19:AT:62:ALA:O	19:AT:63:LYS:HG3	2.05	0.55
23:BB:1168:G:H2'	23:BB:1169:A:O4'	2.06	0.55
23:BB:1241:A:C2'	23:BB:1242:U:H5'	2.36	0.55
23:BB:2053:G:O2'	23:BB:2054:A:H5'	2.06	0.55
23:BB:419:U:H2'	23:BB:420:C:C6	2.41	0.55
23:BB:543:G:C3'	23:BB:544:C:H5''	2.36	0.55
28:BF:121:PHE:HE2	28:BF:166:ARG:H	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:9:GLN:HB2	35:BN:17:ARG:CZ	2.36	0.55
38:BQ:3:VAL:HG12	38:BQ:5:ARG:HG3	1.87	0.55
39:BR:5:PHE:HA	39:BR:40:MET:HG3	1.86	0.55
39:BR:4:VAL:HG22	39:BR:5:PHE:H	1.71	0.55
24:BV:21:ARG:NH2	24:BV:27:PRO:HG3	2.21	0.55
1:CA:1014:A:C2	1:CA:1219:A:H1'	2.41	0.55
1:CA:301:G:H2'	1:CA:302:G:H8	1.71	0.55
10:CK:13:LYS:HD2	10:CK:13:LYS:N	2.21	0.55
23:DB:1460:U:H5''	23:DB:1461:C:C5	2.41	0.55
23:DB:1938:A:O2'	23:DB:1939:U:H5''	2.06	0.55
23:DB:245:G:O2'	23:DB:246:C:H5'	2.07	0.55
23:DB:2538:C:O2'	23:DB:2539:C:H5'	2.06	0.55
25:DC:153:LEU:H	25:DC:153:LEU:HD23	1.71	0.55
26:DD:11:MET:O	26:DD:22:ILE:HA	2.06	0.55
27:DE:48:THR:HG23	27:DE:84:THR:HA	1.88	0.55
52:DI:2:LYS:NZ	52:DI:2:LYS:HB3	2.21	0.55
32:DK:64:ARG:N	32:DK:83:ALA:HB3	2.22	0.55
35:DN:26:GLY:HA2	35:DN:75:ILE:HD13	1.87	0.55
36:DO:49:VAL:HG22	36:DO:50:ALA:N	2.21	0.55
39:DR:42:ALA:HB1	39:DR:53:PHE:CD1	2.40	0.55
1:AA:1202:U:H2'	1:AA:1203:C:H5'	1.87	0.55
1:AA:1397:C:H4'	1:AA:1398:A:OP2	2.04	0.55
1:AA:196:A:O5'	19:AT:63:LYS:HE2	2.05	0.55
1:AA:235:C:H2'	1:AA:236:A:H8	1.70	0.55
1:AA:720:C:H5''	17:AR:40:PRO:HA	1.87	0.55
1:AA:983:A:H5'	1:AA:984:C:OP2	2.06	0.55
1:AA:9:G:H5'	4:AE:107:GLY:HA3	1.86	0.55
8:AI:30:ASN:HD21	8:AI:65:THR:HA	1.71	0.55
10:AK:85:VAL:HG21	10:AK:96:ILE:HD11	1.88	0.55
11:AL:48:LEU:CD2	11:AL:48:LEU:H	2.13	0.55
12:AM:11:HIS:O	12:AM:12:LYS:HG2	2.06	0.55
18:AS:32:THR:HB	18:AS:49:ALA:O	2.05	0.55
19:AT:74:HIS:O	19:AT:78:LEU:HD12	2.05	0.55
23:BB:1244:A:O2'	23:BB:1245:G:H5'	2.05	0.55
23:BB:2398:U:H2'	23:BB:2399:G:H8	1.71	0.55
23:BB:2538:C:O2'	23:BB:2539:C:H5'	2.06	0.55
23:BB:2600:A:O2'	23:BB:2601:C:H5'	2.05	0.55
23:BB:533:G:H2'	23:BB:534:U:C6	2.40	0.55
25:BC:130:PRO:HB2	25:BC:133:ASN:ND2	2.21	0.55
25:BC:209:ALA:HA	25:BC:213:ARG:NE	2.21	0.55
26:BD:122:VAL:HG22	26:BD:128:ARG:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:442:G:H22	27:BE:46:GLN:NE2	2.03	0.55
29:BG:119:GLY:C	29:BG:120:ILE:HD12	2.27	0.55
29:BG:71:LEU:HA	29:BG:74:MET:CG	2.36	0.55
37:BP:83:ILE:HB	37:BP:85:VAL:HG23	1.88	0.55
37:BP:88:ARG:HA	37:BP:88:ARG:HE	1.70	0.55
39:BR:53:PHE:CG	39:BR:54:VAL:N	2.73	0.55
1:CA:1226:C:H4'	1:CA:1227:A:OP1	2.05	0.55
1:CA:922:G:N3	1:CA:1398:A:H2	2.05	0.55
1:CA:1509:C:O2'	1:CA:1510:C:H5'	2.05	0.55
1:CA:21:G:H2'	1:CA:22:G:C8	2.40	0.55
1:CA:123:U:OP1	1:CA:312:C:H5'	2.07	0.55
3:CD:55:ARG:HG3	3:CD:55:ARG:HH11	1.71	0.55
1:CA:1348:U:H4'	8:CI:121:ARG:HG3	1.88	0.55
10:CK:81:LEU:HD23	10:CK:99:LEU:HD21	1.87	0.55
13:CN:56:PRO:O	13:CN:59:GLN:HG3	2.05	0.55
50:D3:12:ARG:CB	50:D3:23:HIS:HB2	2.31	0.55
23:DB:1559:U:H3'	23:DB:1560:G:H5'	1.88	0.55
23:DB:1657:U:OP2	26:DD:141:ARG:HG3	2.07	0.55
23:DB:1923:U:H2'	23:DB:1924:C:C6	2.41	0.55
23:DB:2467:C:O4'	34:DM:118:LYS:HD2	2.07	0.55
23:DB:2586:U:H2'	23:DB:2587:A:C8	2.41	0.55
23:DB:918:A:C2'	23:DB:919:U:H5'	2.30	0.55
25:DC:67:LYS:HG2	25:DC:149:LYS:O	2.05	0.55
30:DH:80:ILE:N	30:DH:80:ILE:HD12	2.21	0.55
34:DM:114:ARG:O	34:DM:117:PHE:HD1	1.89	0.55
34:DM:41:LEU:HB2	34:DM:95:LEU:HD22	1.87	0.55
37:DP:1:SER:HB2	37:DP:5:LYS:HG2	1.88	0.55
1:AA:1241:G:H2'	1:AA:1242:G:C8	2.38	0.55
1:AA:366:A:O2'	1:AA:367:U:OP1	2.24	0.55
12:AM:15:VAL:CG2	12:AM:33:LEU:HD12	2.35	0.55
1:AA:1308:U:OP1	12:AM:95:PRO:HA	2.07	0.55
15:AP:23:ASP:O	15:AP:26:ASN:HB2	2.06	0.55
49:B2:34:ARG:NE	49:B2:42:LEU:HD13	2.21	0.55
23:BB:1199:U:H2'	23:BB:1200:C:C6	2.41	0.55
23:BB:1291:C:O2'	23:BB:1292:G:H5'	2.07	0.55
23:BB:1381:G:H2'	23:BB:1382:G:H5'	1.86	0.55
23:BB:2143:C:O3'	23:BB:2144:G:H4'	2.06	0.55
23:BB:214:G:H2'	23:BB:215:G:C8	2.41	0.55
23:BB:2400:G:O2'	23:BB:2401:U:H5'	2.07	0.55
23:BB:650:C:H5''	50:B3:22:LYS:HZ3	1.71	0.55
23:BB:729:G:H4'	23:BB:763:G:C5'	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:179:GLU:HG3	25:BC:266:ILE:HG22	1.87	0.55
27:BE:147:LEU:HD13	27:BE:182:ALA:O	2.07	0.55
27:BE:187:VAL:HG23	27:BE:187:VAL:O	2.06	0.55
28:BF:33:ILE:HA	28:BF:155:ILE:HA	1.88	0.55
30:BH:4:ILE:H	30:BH:37:VAL:CB	2.17	0.55
32:BK:47:ILE:HG23	32:BK:48:PRO:HD2	1.87	0.55
33:BL:63:LYS:CG	50:B3:11:LYS:HA	2.37	0.55
34:BM:58:LYS:C	34:BM:60:GLN:H	2.07	0.55
35:BN:34:ILE:HG12	35:BN:34:ILE:O	2.07	0.55
37:BP:5:LYS:C	37:BP:7:LEU:H	2.09	0.55
38:BQ:88:GLU:HA	39:BR:53:PHE:CD2	2.41	0.55
43:BW:56:HIS:CE1	43:BW:58:LEU:HB2	2.41	0.55
44:BX:7:ARG:NH1	44:BX:7:ARG:HA	2.21	0.55
45:BY:6:ILE:HD13	45:BY:35:VAL:O	2.06	0.55
1:CA:1217:C:H2'	1:CA:1218:C:H6	1.71	0.55
1:CA:220:G:O2'	1:CA:221:C:H5'	2.06	0.55
1:CA:449:G:H2'	1:CA:450:G:C8	2.40	0.55
20:CB:209:VAL:O	20:CB:213:LEU:HD12	2.05	0.55
5:CF:86:ARG:HD3	17:CR:63:TYR:O	2.06	0.55
19:CT:73:ARG:HG3	19:CT:74:HIS:N	2.21	0.55
23:DB:1022:G:H8	31:DJ:68:LYS:HE3	1.69	0.55
23:DB:1098:A:C3'	52:DI:3:LYS:HB3	2.36	0.55
23:DB:1099:G:O5'	52:DI:3:LYS:C	2.39	0.55
23:DB:192:C:H2'	23:DB:193:U:H5'	1.88	0.55
23:DB:2103:C:H2'	23:DB:2104:C:O4'	2.07	0.55
25:DC:205:GLY:C	25:DC:206:LYS:HG2	2.27	0.55
26:DD:56:LYS:HB3	26:DD:56:LYS:NZ	2.21	0.55
27:DE:45:ALA:O	27:DE:46:GLN:HB3	2.05	0.55
52:DI:79:LEU:HD11	52:DI:131:THR:OG1	2.06	0.55
34:DM:73:ILE:HG21	34:DM:90:GLU:CD	2.27	0.55
35:DN:66:ALA:O	35:DN:70:THR:HB	2.06	0.55
36:DO:26:LEU:O	36:DO:27:VAL:HG13	2.06	0.55
36:DO:67:ASN:HD22	36:DO:68:LYS:N	2.04	0.55
37:DP:29:VAL:O	37:DP:31:VAL:HG23	2.06	0.55
1:AA:1320:C:H41	18:AS:36:ARG:HG2	1.71	0.55
1:AA:1369:C:H2'	1:AA:1370:G:C8	2.42	0.55
1:AA:34:C:H2'	1:AA:35:G:C8	2.41	0.55
1:AA:462:G:H3'	1:AA:463:U:H5''	1.88	0.55
5:AF:64:VAL:CG1	5:AF:65:GLU:H	2.13	0.55
6:AG:41:ILE:HD11	6:AG:116:ALA:HB2	1.88	0.55
1:AA:981:U:H4'	13:AN:60:ARG:HG3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AN:9:GLU:O	13:AN:13:VAL:HG23	2.07	0.55
23:BB:1113:U:H5''	29:BG:2:ARG:N	2.22	0.55
23:BB:1444:G:H2'	23:BB:1445:G:C8	2.40	0.55
23:BB:1854:A:N6	23:BB:1888:G:H1'	2.21	0.55
23:BB:2657:A:H2'	23:BB:2658:C:O4'	2.07	0.55
23:BB:2729:G:H2'	23:BB:2730:C:H6	1.71	0.55
23:BB:2888:C:H2'	23:BB:2889:C:H6	1.70	0.55
23:BB:295:G:O2'	23:BB:296:U:H5'	2.07	0.55
23:BB:621:A:H2'	23:BB:622:G:O4'	2.06	0.55
23:BB:667:U:H2'	23:BB:668:A:O4'	2.07	0.55
23:BB:78:U:H2'	23:BB:79:C:H6	1.70	0.55
28:BF:15:LEU:HA	28:BF:18:GLU:HB3	1.88	0.55
37:BP:24:THR:HG21	37:BP:111:GLU:HG2	1.88	0.55
40:BS:17:VAL:HG23	40:BS:76:VAL:HG11	1.88	0.55
42:BU:23:LYS:O	42:BU:25:LYS:N	2.40	0.55
23:BB:84:A:O5'	42:BU:91:LYS:HD3	2.07	0.55
42:BU:78:LYS:CE	42:BU:96:LYS:HB2	2.37	0.55
23:BB:397:U:OP1	46:BZ:29:GLY:HA2	2.07	0.55
4:CE:65:LYS:HB3	4:CE:65:LYS:NZ	2.21	0.55
6:CG:91:ARG:HB3	6:CG:92:PRO:HD2	1.88	0.55
9:CJ:36:VAL:HA	9:CJ:76:ILE:CG2	2.37	0.55
17:CR:72:ARG:CZ	17:CR:72:ARG:HB2	2.36	0.55
47:D0:26:SER:HB2	47:D0:38:LEU:CD2	2.37	0.55
48:D1:22:THR:HG21	50:D3:34:LYS:NZ	2.21	0.55
23:DB:1099:G:OP2	52:DI:2:LYS:O	2.23	0.55
23:DB:1197:G:H2'	23:DB:1198:U:C6	2.41	0.55
23:DB:1251:C:O2'	23:DB:1252:G:H3'	2.07	0.55
23:DB:1550:C:H2'	23:DB:1551:A:C8	2.41	0.55
23:DB:1551:A:H3'	23:DB:1552:A:H5''	1.87	0.55
23:DB:2250:G:H21	23:DB:2496:C:H4'	1.71	0.55
23:DB:2398:U:H2'	23:DB:2399:G:H8	1.70	0.55
23:DB:2839:G:H2'	23:DB:2840:C:C6	2.41	0.55
23:DB:78:U:H2'	23:DB:79:C:C6	2.41	0.55
23:DB:935:C:H2'	23:DB:936:A:H8	1.71	0.55
25:DC:33:LEU:CD2	25:DC:34:GLU:HG3	2.25	0.55
26:DD:4:LEU:HD23	26:DD:77:ARG:HD3	1.88	0.55
27:DE:59:PRO:HB2	27:DE:60:TRP:CD1	2.41	0.55
30:DH:26:ALA:C	30:DH:28:ASN:N	2.60	0.55
30:DH:3:VAL:HB	30:DH:37:VAL:HG11	1.87	0.55
34:DM:2:LEU:HB2	34:DM:47:GLU:HG2	1.88	0.55
37:DP:23:ASP:HB2	37:DP:93:LYS:HE2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:25:VAL:C	37:DP:27:VAL:H	2.10	0.55
37:DP:86:LYS:HE3	37:DP:88:ARG:HB2	1.89	0.55
39:DR:63:VAL:CG2	39:DR:64:VAL:H	2.09	0.55
42:DU:28:LEU:HD21	42:DU:31:GLY:HA3	1.89	0.55
43:DW:23:LYS:HD3	43:DW:24:ARG:HD2	1.88	0.55
43:DW:35:ILE:HD12	43:DW:35:ILE:N	2.14	0.55
45:DY:36:GLU:O	45:DY:37:ARG:HG2	2.06	0.55
1:AA:1435:G:H2'	1:AA:1436:U:C6	2.41	0.55
1:AA:19:A:OP1	4:AE:134:ASN:ND2	2.39	0.55
1:AA:335:C:H2'	1:AA:336:A:H8	1.72	0.55
1:AA:87:C:H2'	1:AA:88:U:C4'	2.36	0.55
4:AE:13:LYS:HD2	4:AE:112:ALA:HB1	1.89	0.55
15:AP:1:MET:N	15:AP:24:SER:HB3	2.22	0.55
47:B0:9:ARG:NH1	47:B0:9:ARG:HB3	2.22	0.55
49:B2:35:ARG:O	49:B2:39:ARG:HA	2.06	0.55
23:BB:1360:G:C2'	23:BB:1361:G:H5'	2.37	0.55
23:BB:1564:C:O2'	23:BB:1565:C:H5'	2.06	0.55
23:BB:2639:A:H2'	23:BB:2640:G:O4'	2.06	0.55
23:BB:2849:U:H4'	23:BB:2850:A:C5'	2.37	0.55
23:BB:338:G:C2	23:BB:339:U:H1'	2.41	0.55
23:BB:363:G:H2'	23:BB:364:C:C6	2.42	0.55
23:BB:654:A:H2'	23:BB:655:A:C5'	2.33	0.55
25:BC:254:LYS:N	25:BC:254:LYS:HD2	2.22	0.55
27:BE:48:THR:HG23	27:BE:87:ALA:H	1.72	0.55
28:BF:174:PHE:N	28:BF:175:PRO:HD2	2.22	0.55
29:BG:5:LYS:HB3	29:BG:68:ARG:CZ	2.36	0.55
33:BL:30:THR:HG22	33:BL:36:LYS:HZ1	1.72	0.55
23:BB:2848:G:O2'	37:BP:96:LEU:HD22	2.06	0.55
38:BQ:5:ARG:HH11	38:BQ:5:ARG:CB	2.13	0.55
40:BS:29:VAL:HG11	40:BS:69:LEU:HB2	1.88	0.55
41:BT:35:ALA:O	41:BT:81:LYS:HD3	2.06	0.55
24:BV:2:PHE:O	24:BV:3:THR:HG23	2.07	0.55
46:BZ:33:ASN:CB	46:BZ:44:PHE:HB2	2.29	0.55
1:CA:214:C:H2'	1:CA:215:C:H6	1.72	0.55
1:CA:279:A:H4'	1:CA:280:C:OP2	2.07	0.55
1:CA:399:G:H2'	1:CA:400:C:C6	2.42	0.55
1:CA:806:C:H2'	1:CA:807:A:H8	1.71	0.55
3:CD:44:LYS:O	3:CD:44:LYS:HD3	2.07	0.55
10:CK:19:VAL:CG2	10:CK:34:THR:HG22	2.31	0.55
11:CL:73:LEU:HD21	11:CL:103:CYS:HA	1.89	0.55
1:CA:1526:G:H5''	21:CU:38:GLU:HG2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:D1:22:THR:O	48:D1:23:THR:C	2.44	0.55
51:D4:26:ILE:HD11	51:D4:34:LYS:HA	1.88	0.55
22:DA:51:G:O2'	22:DA:52:A:H5''	2.06	0.55
23:DB:1061:U:H4'	23:DB:1070:A:O3'	2.07	0.55
23:DB:1803:A:H5'	25:DC:255:LYS:HD3	1.88	0.55
23:DB:2262:U:OP1	23:DB:2387:U:O2'	2.23	0.55
23:DB:2756:U:H4'	23:DB:2757:A:OP1	2.06	0.55
23:DB:2825:G:H2'	23:DB:2826:A:H5'	1.89	0.55
23:DB:639:U:H2'	23:DB:640:C:H6	1.69	0.55
23:DB:947:A:O2'	23:DB:984:A:H2	1.90	0.55
25:DC:19:VAL:HG12	25:DC:20:ASN:H	1.70	0.55
26:DD:138:LEU:CD1	26:DD:142:VAL:HB	2.37	0.55
26:DD:18:ASP:C	26:DD:20:VAL:H	2.10	0.55
26:DD:5:VAL:HG12	26:DD:6:GLY:N	2.21	0.55
35:DN:110:MET:O	35:DN:111:ALA:HB3	2.07	0.55
39:DR:66:HIS:HA	39:DR:98:ILE:HA	1.89	0.55
42:DU:11:ILE:HB	42:DU:69:VAL:CG2	2.35	0.55
42:DU:27:VAL:HG12	42:DU:33:VAL:HG13	1.87	0.55
43:DW:18:LYS:HB2	43:DW:18:LYS:HZ2	1.69	0.55
1:AA:913:A:H1'	1:AA:914:A:O4'	2.06	0.55
2:AC:166:TRP:HA	2:AC:166:TRP:CE3	2.41	0.55
16:AQ:63:CYS:SG	16:AQ:73:THR:HG23	2.47	0.55
18:AS:49:ALA:HB1	18:AS:56:HIS:CB	2.31	0.55
23:BB:142:A:O2'	23:BB:143:C:H5'	2.07	0.55
23:BB:1551:A:H3'	23:BB:1552:A:H5''	1.87	0.55
23:BB:1754:A:OP2	37:BP:94:ALA:HB1	2.06	0.55
23:BB:2047:C:H2'	23:BB:2048:G:C8	2.42	0.55
23:BB:2393:U:H5'	33:BL:62:PRO:CD	2.36	0.55
23:BB:2471:A:O2'	23:BB:2472:G:C8	2.54	0.55
23:BB:2557:G:H2'	23:BB:2558:C:H6	1.71	0.55
23:BB:395:U:H2'	23:BB:396:G:N7	2.21	0.55
23:BB:566:U:O2'	23:BB:567:U:H5'	2.06	0.55
26:BD:31:ALA:HA	26:BD:51:THR:OG1	2.07	0.55
26:BD:93:GLY:O	26:BD:94:GLN:HG3	2.06	0.55
29:BG:89:VAL:HG12	29:BG:159:LYS:HA	1.89	0.55
31:BJ:112:GLY:O	31:BJ:116:ARG:HB2	2.06	0.55
33:BL:27:LEU:C	33:BL:29:LYS:H	2.10	0.55
34:BM:7:THR:HG22	34:BM:8:LYS:N	2.17	0.55
36:BO:56:LYS:HE3	36:BO:57:ALA:N	2.22	0.55
40:BS:70:LYS:HD2	40:BS:110:ARG:HG3	1.88	0.55
42:BU:5:ARG:HG2	42:BU:6:ARG:N	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BW:45:HIS:HB2	43:BW:67:LYS:CE	2.37	0.55
1:CA:1221:G:O3'	18:CS:76:THR:HG21	2.06	0.55
1:CA:1226:C:O2'	1:CA:1227:A:H5'	2.07	0.55
1:CA:129:A:H1'	1:CA:130:A:C8	2.41	0.55
1:CA:1335:U:H5''	1:CA:1337:G:N2	2.21	0.55
1:CA:219:U:H2'	1:CA:220:G:C8	2.42	0.55
1:CA:389:A:H2'	1:CA:389:A:N3	2.22	0.55
1:CA:462:G:H3'	1:CA:463:U:H5''	1.88	0.55
1:CA:824:G:H2'	1:CA:825:A:H8	1.71	0.55
3:CD:117:VAL:HG12	3:CD:130:ASN:CA	2.37	0.55
1:CA:412:A:N1	3:CD:30:LYS:HG3	2.20	0.55
5:CF:81:ASN:O	5:CF:84:VAL:HG12	2.06	0.55
8:CI:49:GLN:N	8:CI:50:PRO:HD2	2.22	0.55
12:CM:44:ILE:HA	12:CM:47:LEU:HD21	1.89	0.55
16:CQ:6:THR:HG23	16:CQ:59:GLU:OE1	2.06	0.55
18:CS:11:ASP:HB3	18:CS:13:HIS:CE1	2.42	0.55
47:D0:41:HIS:CG	47:D0:42:ILE:N	2.70	0.55
50:D3:2:LYS:HB2	50:D3:2:LYS:HZ2	1.72	0.55
23:DB:1099:G:P	52:DI:4:VAL:N	2.65	0.55
23:DB:1219:U:H2'	23:DB:1220:G:C8	2.41	0.55
23:DB:1405:U:H2'	23:DB:1406:U:H6	1.72	0.55
23:DB:2078:C:O2'	23:DB:2079:U:H5'	2.06	0.55
23:DB:2230:G:H2'	23:DB:2231:U:C6	2.42	0.55
23:DB:2642:G:OP1	31:DJ:84:ILE:HG21	2.06	0.55
23:DB:705:A:N6	23:DB:726:G:O2'	2.39	0.55
23:DB:1824:G:O2'	25:DC:244:VAL:HG21	2.06	0.55
27:DE:149:ILE:HD11	27:DE:187:VAL:H	1.72	0.55
29:DG:10:VAL:HG12	29:DG:10:VAL:O	2.05	0.55
31:DJ:135:GLN:NE2	31:DJ:137:PRO:HB2	2.19	0.55
31:DJ:44:TYR:HE2	31:DJ:50:THR:HB	1.71	0.55
33:DL:108:ALA:HB3	33:DL:125:LEU:CB	2.37	0.55
34:DM:133:LYS:HD2	34:DM:134:THR:H	1.71	0.55
34:DM:71:LYS:O	34:DM:73:ILE:HG12	2.07	0.55
38:DQ:23:TYR:HB2	38:DQ:28:SER:HB3	1.89	0.55
40:DS:34:ASP:HA	40:DS:37:THR:HG22	1.88	0.55
42:DU:94:PHE:CD2	42:DU:100:GLU:HG2	2.42	0.55
24:DV:8:VAL:HG12	24:DV:38:LEU:HD11	1.89	0.55
44:DX:23:ARG:HA	44:DX:26:PHE:CD1	2.42	0.55
23:DB:2080:A:H5'	46:DZ:17:SER:CB	2.37	0.55
1:AA:1289:A:C2'	1:AA:1290:G:H5'	2.35	0.55
1:AA:173:U:H5'	1:AA:197:A:O4'	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:220:G:O2'	1:AA:221:C:H5'	2.06	0.55
1:AA:70:U:H1'	1:AA:71:A:N7	2.22	0.55
1:AA:801:U:H2'	1:AA:802:A:C8	2.42	0.55
1:AA:806:C:H2'	1:AA:807:A:H8	1.72	0.55
3:AD:152:SER:O	3:AD:155:LYS:HG3	2.07	0.55
4:AE:80:LEU:HB2	4:AE:97:PRO:HB3	1.87	0.55
6:AG:71:THR:O	6:AG:72:VAL:HG13	2.06	0.55
8:AI:56:MET:SD	8:AI:57:VAL:N	2.80	0.55
16:AQ:58:VAL:HB	16:AQ:74:LEU:HD23	1.88	0.55
22:BA:6:G:H2'	22:BA:7:G:H8	1.69	0.55
23:BB:1515:A:H5'	23:BB:1557:C:H5'	1.88	0.55
23:BB:2138:G:H2'	23:BB:2139:U:C6	2.42	0.55
23:BB:36:G:N2	23:BB:37:C:H1'	2.22	0.55
23:BB:458:G:P	49:B2:39:ARG:HH22	2.29	0.55
23:BB:589:U:H2'	23:BB:590:A:H8	1.72	0.55
23:BB:921:C:O2'	23:BB:922:C:H5'	2.06	0.55
27:BE:118:LEU:C	27:BE:119:ILE:HD12	2.27	0.55
29:BG:77:GLY:HA3	29:BG:135:ALA:O	2.07	0.55
30:BH:75:LEU:HD23	30:BH:76:GLU:H	1.71	0.55
32:BK:97:THR:OG1	32:BK:98:ARG:HD2	2.06	0.55
36:BO:76:LYS:HB2	36:BO:106:LEU:HD22	1.88	0.55
24:BV:77:VAL:HG12	34:BM:133:LYS:NZ	2.21	0.55
1:CA:1157:A:H5'	1:CA:1158:C:C6	2.42	0.55
1:CA:230:G:O2'	1:CA:231:U:H5'	2.07	0.55
6:CG:106:ALA:HB1	6:CG:132:THR:OG1	2.07	0.55
8:CI:50:PRO:O	8:CI:54:VAL:HG22	2.06	0.55
10:CK:17:ASP:OD1	10:CK:36:ARG:HG3	2.06	0.55
11:CL:54:VAL:CG1	11:CL:79:ILE:HD11	2.37	0.55
18:CS:30:LEU:HD12	18:CS:48:ILE:HG12	1.88	0.55
47:D0:41:HIS:NE2	47:D0:42:ILE:HG22	2.22	0.55
47:D0:42:ILE:O	47:D0:46:GLY:N	2.39	0.55
23:DB:1444:G:H2'	23:DB:1445:G:C8	2.40	0.55
23:DB:2213:U:O2	23:DB:2213:U:H2'	2.07	0.55
23:DB:2282:G:O2'	23:DB:2283:C:P	2.64	0.55
23:DB:278:A:N3	23:DB:278:A:H2'	2.20	0.55
23:DB:321:U:OP2	27:DE:130:LYS:HA	2.07	0.55
23:DB:667:U:H2'	23:DB:668:A:O4'	2.07	0.55
23:DB:845:A:C6	23:DB:847:U:H1'	2.42	0.55
23:DB:90:U:OP2	23:DB:91:A:H3'	2.07	0.55
25:DC:216:ARG:HB3	25:DC:217:PRO:CD	2.37	0.55
26:DD:118:PHE:HA	26:DD:164:GLN:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DF:98:PHE:HA	28:DF:101:ARG:NE	2.18	0.55
52:DI:92:PRO:O	52:DI:93:ASN:HB2	2.07	0.55
23:DB:1651:G:H4'	35:DN:39:PRO:HG2	1.89	0.55
35:DN:4:ARG:NH2	35:DN:4:ARG:H	2.05	0.55
37:DP:52:ARG:HD2	37:DP:60:VAL:HG13	1.88	0.55
40:DS:72:THR:HB	40:DS:109:ASP:HB2	1.88	0.55
1:AA:219:U:H2'	1:AA:220:G:C8	2.42	0.55
1:AA:279:A:H4'	1:AA:280:C:OP2	2.06	0.55
1:AA:389:A:H2'	1:AA:389:A:N3	2.22	0.55
1:AA:426:U:H2'	1:AA:427:U:C6	2.41	0.55
20:AB:30:ILE:HA	20:AB:39:ILE:O	2.07	0.55
3:AD:115:GLN:HG3	3:AD:119:HIS:CE1	2.41	0.55
8:AI:56:MET:HA	8:AI:59:LYS:NZ	2.22	0.55
10:AK:15:VAL:HG11	10:AK:35:ASP:HB2	1.88	0.55
1:AA:521:G:OP2	11:AL:50:LYS:HE3	2.07	0.55
14:AO:78:THR:HA	14:AO:81:ILE:HG12	1.89	0.55
16:AQ:18:LYS:HA	16:AQ:47:ASP:O	2.07	0.55
17:AR:21:ASP:OD2	17:AR:23:LYS:HG3	2.06	0.55
48:B1:31:GLU:CG	48:B1:32:LYS:H	2.19	0.55
23:BB:1405:U:H2'	23:BB:1406:U:H6	1.71	0.55
23:BB:1591:A:H2'	23:BB:1592:C:C6	2.42	0.55
23:BB:1904:G:O2'	23:BB:1905:C:H5'	2.06	0.55
23:BB:2236:U:H2'	23:BB:2237:G:O4'	2.07	0.55
23:BB:2265:U:H3'	23:BB:2266:A:C5'	2.36	0.55
23:BB:2307:G:H4'	23:BB:2311:A:N6	2.22	0.55
23:BB:280:U:H2'	23:BB:281:C:C6	2.42	0.55
23:BB:2839:G:H2'	23:BB:2840:C:C6	2.41	0.55
23:BB:329:G:O6	42:BU:16:LYS:HE2	2.07	0.55
23:BB:856:G:H21	43:BW:22:VAL:HG11	1.72	0.55
25:BC:42:ARG:CG	25:BC:43:ASN:N	2.62	0.55
26:BD:48:ILE:HG13	26:BD:80:TRP:CD1	2.41	0.55
52:BI:49:GLU:CG	52:BI:54:ILE:HD11	2.37	0.55
32:BK:87:LEU:HA	32:BK:95:ILE:H	1.72	0.55
23:BB:635:C:H3'	33:BL:109:LYS:HZ1	1.70	0.55
33:BL:77:ILE:CB	33:BL:110:VAL:HG22	2.31	0.55
33:BL:63:LYS:HB2	50:B3:11:LYS:HG3	1.87	0.55
34:BM:42:THR:O	34:BM:44:ARG:N	2.39	0.55
23:BB:2335:A:OP2	36:BO:9:ARG:HG3	2.07	0.55
39:BR:22:LEU:O	39:BR:96:VAL:HG22	2.06	0.55
41:BT:68:LYS:C	41:BT:68:LYS:HD3	2.28	0.55
1:CA:1062:U:H2'	1:CA:1063:C:C6	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1386:G:H2'	1:CA:1387:G:C8	2.42	0.55
1:CA:16:A:N1	1:CA:919:A:H2	2.04	0.55
1:CA:254:G:O2'	1:CA:255:G:H5'	2.07	0.55
20:CB:205:ALA:HB3	20:CB:208:ALA:HB3	1.88	0.55
3:CD:78:ALA:HB1	3:CD:85:THR:HA	1.88	0.55
1:CA:1248:A:H2	8:CI:71:ILE:HD11	1.70	0.55
11:CL:120:ARG:HD3	11:CL:121:PRO:HD2	1.88	0.55
12:CM:44:ILE:C	12:CM:44:ILE:HD12	2.27	0.55
13:CN:68:ARG:NH2	13:CN:80:ARG:HH12	2.05	0.55
23:DB:1432:G:O2'	23:DB:1433:A:H5'	2.07	0.55
23:DB:1564:C:O2'	23:DB:1565:C:H5'	2.06	0.55
23:DB:2216:G:H2'	23:DB:2217:G:H8	1.72	0.55
23:DB:2895:G:H2'	23:DB:2896:C:H6	1.66	0.55
23:DB:741:U:H2'	23:DB:742:A:C8	2.41	0.55
27:DE:49:ARG:C	27:DE:51:GLU:H	2.09	0.55
29:DG:79:THR:HG23	29:DG:80:GLU:N	2.22	0.55
32:DK:108:ARG:NH2	37:DP:36:LYS:H	2.04	0.55
37:DP:59:THR:HA	37:DP:76:HIS:HA	1.89	0.55
37:DP:93:LYS:CG	37:DP:96:LEU:HA	2.26	0.55
39:DR:41:ILE:O	39:DR:43:ASN:N	2.39	0.55
43:DW:20:LEU:HD11	43:DW:31:LEU:HD12	1.88	0.55
46:DZ:3:LYS:HD2	46:DZ:3:LYS:H	1.72	0.55
1:AA:1307:U:H2'	1:AA:1308:U:C6	2.42	0.55
1:AA:415:A:H3'	1:AA:416:G:H8	1.72	0.55
20:AB:53:LEU:HD11	20:AB:216:VAL:HA	1.87	0.55
3:AD:98:ASP:HB2	3:AD:132:ALA:HB1	1.88	0.55
6:AG:99:ALA:O	6:AG:103:ILE:HG13	2.07	0.55
8:AI:64:ILE:HD13	8:AI:78:ILE:CG2	2.37	0.55
13:AN:61:ASN:O	13:AN:62:ARG:HB2	2.07	0.55
23:BB:1327:A:H2'	23:BB:1328:A:O4'	2.06	0.55
23:BB:2093:G:O2'	23:BB:2094:A:H5'	2.07	0.55
23:BB:235:U:H2'	23:BB:236:C:C6	2.42	0.55
23:BB:62:U:H3'	23:BB:63:A:H8	1.72	0.55
23:BB:64:A:H2'	23:BB:65:U:C6	2.42	0.55
23:BB:866:A:H61	23:BB:913:U:C4'	2.19	0.55
23:BB:1843:C:H5''	25:BC:252:LYS:HZ3	1.72	0.55
25:BC:260:LYS:HB2	25:BC:261:ARG:HE	1.71	0.55
28:BF:165:GLY:HA2	28:BF:168:LEU:HD21	1.88	0.55
30:BH:11:ASN:ND2	30:BH:12:LEU:N	2.55	0.55
23:BB:1131:G:C4'	31:BJ:85:LYS:HZ2	2.20	0.55
37:BP:112:ARG:HH11	37:BP:112:ARG:CA	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:1:SER:HA	37:BP:4:ILE:HD12	1.88	0.55
37:BP:23:ASP:N	37:BP:93:LYS:HE2	2.22	0.55
42:BU:11:ILE:HG23	42:BU:20:LYS:O	2.07	0.55
43:BW:21:GLY:HA3	43:BW:32:ALA:CB	2.36	0.55
1:CA:1068:G:N7	1:CA:1094:G:H2'	2.22	0.55
1:CA:118:U:O4	1:CA:288:A:H2'	2.07	0.55
1:CA:454:G:O2'	1:CA:455:G:H5'	2.07	0.55
1:CA:985:C:H2'	1:CA:986:U:H6	1.71	0.55
3:CD:118:SER:HA	3:CD:130:ASN:HB2	1.87	0.55
9:CJ:40:ILE:HG13	9:CJ:73:LEU:HB3	1.89	0.55
11:CL:17:LYS:HZ2	11:CL:17:LYS:HB2	1.71	0.55
16:CQ:23:ALA:C	16:CQ:24:ILE:HD12	2.27	0.55
16:CQ:68:LYS:O	16:CQ:70:LYS:N	2.40	0.55
16:CQ:74:LEU:HD22	16:CQ:75:VAL:N	2.22	0.55
22:DA:103:U:O2'	22:DA:104:A:H5'	2.07	0.55
22:DA:76:G:H2'	22:DA:77:U:C6	2.38	0.55
23:DB:1351:C:O2'	23:DB:1571:A:H1'	2.06	0.55
23:DB:1921:G:O2'	23:DB:1922:G:H5'	2.07	0.55
23:DB:2400:G:O2'	23:DB:2401:U:H5'	2.07	0.55
23:DB:2848:G:H1'	23:DB:2868:A:H61	1.72	0.55
23:DB:2239:G:OP1	25:DC:246:PRO:HG3	2.07	0.55
27:DE:112:LEU:HD12	27:DE:115:GLN:NE2	2.22	0.55
29:DG:34:ARG:NH1	29:DG:70:LEU:HG	2.22	0.55
30:DH:4:ILE:HD13	30:DH:4:ILE:H	1.71	0.55
31:DJ:44:TYR:CD1	31:DJ:45:THR:N	2.71	0.55
39:DR:86:GLN:HG3	39:DR:87:GLN:HG2	1.89	0.55
40:DS:58:ALA:O	40:DS:62:ASP:HB3	2.06	0.55
40:DS:85:ILE:HD13	40:DS:86:MET:N	2.21	0.55
41:DT:48:GLN:HG3	41:DT:49:LYS:N	2.21	0.55
41:DT:53:VAL:HA	41:DT:93:LEU:HG	1.88	0.55
42:DU:27:VAL:HG12	42:DU:33:VAL:HG22	1.87	0.55
42:DU:96:LYS:HD3	42:DU:97:SER:N	2.21	0.55
24:DV:26:PHE:HE2	24:DV:44:HIS:HA	1.72	0.55
43:DW:47:GLY:CA	43:DW:71:LYS:HB3	2.34	0.55
1:AA:1033:G:H2'	1:AA:1034:G:O4'	2.07	0.55
1:AA:1494:G:O2'	1:AA:1495:U:H5'	2.06	0.55
1:AA:921:U:H2'	1:AA:922:G:O4'	2.07	0.55
7:AH:31:LEU:O	7:AH:35:ILE:HG13	2.07	0.55
15:AP:18:GLN:NE2	15:AP:35:ARG:HD2	2.22	0.55
16:AQ:20:ILE:HG21	16:AQ:52:CYS:SG	2.47	0.55
1:AA:132:C:H5''	19:AT:68:LYS:CE	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:B1:10:LEU:HD13	48:B1:25:ASN:HD22	1.71	0.55
22:BA:91:C:O2'	22:BA:92:C:H5'	2.07	0.55
23:BB:1001:A:H2'	23:BB:1002:G:O4'	2.07	0.55
23:BB:1300:G:H5'	23:BB:1301:A:N3	2.22	0.55
23:BB:1439:A:C6	23:BB:1552:A:C5	2.95	0.55
23:BB:2104:C:H2'	23:BB:2105:U:H5	1.72	0.55
23:BB:607:U:O4	23:BB:620:G:H5''	2.07	0.55
25:BC:68:ARG:HH22	25:BC:190:THR:HG23	1.72	0.55
25:BC:80:LEU:HA	25:BC:91:ALA:HA	1.89	0.55
26:BD:35:THR:CG2	26:BD:46:ARG:HH22	2.20	0.55
28:BF:35:LEU:HD22	28:BF:56:LEU:HD21	1.88	0.55
30:BH:121:VAL:HG21	30:BH:128:HIS:HE2	1.72	0.55
33:BL:39:LYS:CB	33:BL:46:VAL:HG13	2.37	0.55
33:BL:77:ILE:HD12	33:BL:110:VAL:CG2	2.37	0.55
33:BL:77:ILE:CD1	33:BL:92:LEU:HD13	2.32	0.55
36:BO:12:THR:OG1	36:BO:13:ARG:HD2	2.07	0.55
23:BB:994:C:H3'	38:BQ:53:LYS:HE2	1.88	0.55
44:BX:38:GLN:O	44:BX:39:GLN:HG3	2.07	0.55
45:BY:6:ILE:HD13	45:BY:6:ILE:N	2.22	0.55
1:CA:1297:G:H4'	1:CA:1298:U:H5'	1.88	0.55
1:CA:1355:G:H2'	1:CA:1356:G:C8	2.41	0.55
1:CA:610:U:O4'	1:CA:610:U:O2	2.25	0.55
20:CB:101:THR:HG22	20:CB:174:GLU:OE1	2.06	0.55
2:CC:58:ARG:HA	2:CC:63:ILE:HA	1.88	0.55
3:CD:16:THR:HG22	3:CD:17:ASP:H	1.71	0.55
6:CG:91:ARG:O	6:CG:95:ARG:HG3	2.07	0.55
10:CK:106:ILE:HD11	10:CK:109:ILE:HG13	1.87	0.55
1:CA:500:G:H5''	11:CL:120:ARG:HH21	1.71	0.55
12:CM:80:MET:HA	12:CM:87:GLY:CA	2.37	0.55
13:CN:82:LYS:O	13:CN:85:GLU:HB2	2.07	0.55
19:CT:66:ILE:HG23	19:CT:70:LYS:HG3	1.87	0.55
22:DA:66:A:O2'	22:DA:67:G:H8	1.90	0.55
23:DB:1059:G:H4'	52:DI:116:MET:HE2	1.89	0.55
23:DB:106:C:H2'	23:DB:107:G:H8	1.71	0.55
23:DB:1300:G:H5'	23:DB:1301:A:N3	2.22	0.55
23:DB:1934:C:O2'	23:DB:1935:G:H5'	2.07	0.55
23:DB:2228:G:H2'	23:DB:2229:U:C6	2.42	0.55
23:DB:2553:G:H2'	23:DB:2554:U:H4'	1.89	0.55
23:DB:2729:G:H2'	23:DB:2730:C:H6	1.72	0.55
23:DB:2843:G:O2'	23:DB:2844:G:H5'	2.07	0.55
29:DG:126:THR:HG23	29:DG:129:GLU:H	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:DH:5:LEU:HD13	30:DH:35:LYS:CE	2.37	0.55
52:DI:102:ARG:HG3	52:DI:141:ASP:HB2	1.89	0.55
31:DJ:41:LYS:NZ	31:DJ:45:THR:HA	2.22	0.55
32:DK:64:ARG:HD3	32:DK:83:ALA:HB2	1.89	0.55
33:DL:109:LYS:O	33:DL:110:VAL:HG13	2.07	0.55
40:DS:86:MET:HG3	40:DS:87:PRO:HD2	1.89	0.55
41:DT:31:VAL:HG13	41:DT:32:LEU:N	2.16	0.55
24:DV:24:ASN:HB3	24:DV:45:ASP:OD1	2.06	0.55
1:AA:1110:A:H2'	1:AA:1111:A:H5'	1.89	0.54
1:AA:118:U:O4	1:AA:288:A:H2'	2.07	0.54
3:AD:111:ALA:O	3:AD:114:ARG:HB3	2.07	0.54
12:AM:3:ILE:HG23	12:AM:56:ARG:HG2	1.87	0.54
13:AN:33:VAL:HG23	13:AN:40:ARG:NH2	2.23	0.54
23:BB:1166:G:H2'	23:BB:1167:C:C6	2.43	0.54
23:BB:1565:C:OP1	25:BC:23:LEU:HD23	2.08	0.54
23:BB:1637:A:H2'	23:BB:1638:C:H6	1.73	0.54
23:BB:1774:C:C2'	23:BB:1774:C:O2	2.54	0.54
23:BB:1921:G:O2'	23:BB:1922:G:H5'	2.07	0.54
23:BB:528:A:C2	23:BB:2043:C:H4'	2.42	0.54
23:BB:2320:U:H3'	23:BB:2321:U:H5'	1.88	0.54
23:BB:372:G:N2	23:BB:401:A:OP2	2.35	0.54
23:BB:26:G:H1'	23:BB:514:A:H61	1.70	0.54
23:BB:593:U:H2'	23:BB:594:U:C6	2.41	0.54
23:BB:729:G:H4'	23:BB:763:G:H5'	1.88	0.54
23:BB:764:A:H5''	25:BC:208:GLY:CA	2.37	0.54
23:BB:1821:A:H5'	25:BC:155:ARG:HD2	1.89	0.54
23:BB:2575:C:H5''	26:BD:149:ASN:N	2.22	0.54
26:BD:167:ASN:N	26:BD:167:ASN:HD22	2.04	0.54
27:BE:49:ARG:HB2	27:BE:49:ARG:NH1	2.23	0.54
28:BF:37:MET:HG2	28:BF:86:CYS:HB2	1.89	0.54
30:BH:69:ALA:HB2	30:BH:139:PHE:O	2.07	0.54
52:BI:74:PRO:O	52:BI:77:VAL:HG22	2.07	0.54
31:BJ:37:ARG:NH2	31:BJ:110:PRO:HG3	2.22	0.54
33:BL:79:LEU:CA	33:BL:113:ALA:HB3	2.37	0.54
44:BX:38:GLN:H	44:BX:38:GLN:HE21	1.54	0.54
1:CA:1077:G:N2	1:CA:1080:A:OP2	2.40	0.54
1:CA:1320:C:H41	18:CS:36:ARG:HG3	1.72	0.54
4:CE:108:GLY:O	4:CE:111:ARG:HB3	2.06	0.54
6:CG:134:VAL:O	6:CG:138:GLU:HG3	2.06	0.54
1:CA:598:U:H4'	7:CH:85:TYR:CD2	2.43	0.54
14:CO:80:LEU:O	14:CO:84:LEU:HD13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CP:28:ARG:HD3	15:CP:29:ASN:OD1	2.07	0.54
16:CQ:83:LEU:HD22	16:CQ:83:LEU:N	2.22	0.54
17:CR:54:LEU:O	17:CR:58:ILE:HG13	2.07	0.54
21:CU:24:LYS:HB3	21:CU:24:LYS:NZ	2.22	0.54
48:D1:5:ARG:HA	48:D1:28:THR:OG1	2.07	0.54
50:D3:5:THR:O	50:D3:5:THR:HG23	2.06	0.54
22:DA:49:C:H2'	22:DA:50:A:C8	2.42	0.54
23:DB:1022:G:N2	23:DB:1142:A:C2	2.75	0.54
23:DB:1291:C:O2'	23:DB:1292:G:H5'	2.07	0.54
23:DB:1400:U:H2'	23:DB:1401:G:C8	2.41	0.54
23:DB:1773:A:N6	25:DC:206:LYS:HE2	2.22	0.54
23:DB:191:A:O2'	23:DB:192:C:H5'	2.08	0.54
23:DB:2097:A:H2'	23:DB:2098:U:C6	2.42	0.54
23:DB:2646:C:H2'	23:DB:2647:U:O4'	2.07	0.54
23:DB:810:U:C2	33:DL:37:GLY:HA2	2.43	0.54
25:DC:37:SER:H	25:DC:62:ARG:HB3	1.71	0.54
27:DE:164:LEU:HD22	27:DE:164:LEU:O	2.06	0.54
28:DF:140:ILE:N	28:DF:140:ILE:HD12	2.17	0.54
29:DG:123:GLU:HG2	29:DG:131:VAL:HG13	1.89	0.54
30:DH:113:SER:HB2	30:DH:132:PHE:CZ	2.42	0.54
52:DI:49:GLU:CB	52:DI:52:LEU:HD12	2.37	0.54
52:DI:99:LYS:HD3	52:DI:99:LYS:H	1.71	0.54
31:DJ:120:ARG:HB3	31:DJ:121:LYS:NZ	2.22	0.54
31:DJ:120:ARG:N	31:DJ:121:LYS:HZ2	2.05	0.54
32:DK:43:ILE:H	32:DK:43:ILE:CD1	2.19	0.54
33:DL:123:ARG:HA	33:DL:142:ILE:HA	1.88	0.54
23:DB:635:C:H3'	33:DL:126:ARG:HH21	1.73	0.54
34:DM:101:VAL:HG12	34:DM:102:LEU:N	2.21	0.54
35:DN:4:ARG:H	35:DN:4:ARG:CZ	2.19	0.54
37:DP:47:ILE:HD13	37:DP:63:ILE:HG12	1.89	0.54
41:DT:34:VAL:HG22	41:DT:35:ALA:N	2.20	0.54
41:DT:92:ASN:O	41:DT:93:LEU:HD12	2.06	0.54
43:DW:36:ILE:HG21	43:DW:68:PHE:CE1	2.42	0.54
1:AA:944:G:H1'	1:AA:1339:A:N6	2.23	0.54
1:AA:254:G:O2'	1:AA:255:G:H5'	2.07	0.54
1:AA:123:U:OP1	1:AA:312:C:H5'	2.07	0.54
1:AA:915:A:H2'	1:AA:916:U:H5'	1.89	0.54
1:AA:98:A:H2'	1:AA:99:C:C6	2.42	0.54
5:AF:29:ILE:HD13	5:AF:64:VAL:HG11	1.88	0.54
10:AK:109:ILE:CG2	21:AU:16:ARG:HD2	2.37	0.54
12:AM:95:PRO:HD3	12:AM:108:ARG:HG2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AP:52:LEU:HG	15:AP:75:ILE:HG12	1.89	0.54
51:B4:25:VAL:HB	51:B4:33:HIS:C	2.26	0.54
22:BA:76:G:H2'	22:BA:77:U:C6	2.42	0.54
23:BB:1028:A:H2'	23:BB:1029:A:C8	2.42	0.54
23:BB:1227:G:O2'	23:BB:1228:G:H5'	2.08	0.54
23:BB:2292:U:H2'	23:BB:2293:G:H8	1.73	0.54
23:BB:2467:C:C1'	34:BM:118:LYS:HG2	2.37	0.54
23:BB:2512:C:H2'	23:BB:2513:A:O4'	2.07	0.54
23:BB:2646:C:H2'	23:BB:2647:U:O4'	2.07	0.54
25:BC:70:LYS:HE3	25:BC:95:TYR:CD2	2.43	0.54
23:BB:2053:G:OP1	26:BD:149:ASN:C	2.44	0.54
27:BE:143:LEU:CG	27:BE:185:LYS:HD2	2.33	0.54
28:BF:52:ALA:HA	28:BF:148:VAL:HG13	1.90	0.54
28:BF:7:TYR:HA	28:BF:11:VAL:HB	1.90	0.54
29:BG:154:GLU:OE1	29:BG:156:TYR:HB2	2.07	0.54
32:BK:98:ARG:HA	32:BK:118:LEU:HD22	1.89	0.54
34:BM:40:ARG:HD3	34:BM:92:TRP:HB3	1.90	0.54
38:BQ:52:ARG:NH1	38:BQ:55:GLN:HE21	2.05	0.54
23:BB:559:G:H1'	38:BQ:55:GLN:HE22	1.72	0.54
39:BR:39:LEU:O	39:BR:54:VAL:HG12	2.06	0.54
40:BS:44:ALA:O	40:BS:47:VAL:HG22	2.07	0.54
42:BU:33:VAL:HG21	42:BU:66:VAL:HA	1.89	0.54
42:BU:13:LEU:HG	42:BU:69:VAL:HG22	1.89	0.54
1:CA:1064:G:H4'	1:CA:1065:U:OP1	2.06	0.54
1:CA:1217:C:H2'	1:CA:1218:C:C6	2.41	0.54
1:CA:1476:A:H2'	1:CA:1477:U:O4'	2.07	0.54
1:CA:801:U:H2'	1:CA:802:A:C8	2.42	0.54
20:CB:76:SER:O	20:CB:80:LYS:HG2	2.06	0.54
5:CF:92:THR:HG22	5:CF:93:LYS:H	1.73	0.54
5:CF:86:ARG:HH11	17:CR:63:TYR:HB3	1.72	0.54
47:D0:52:LYS:HE2	47:D0:53:VAL:HB	1.89	0.54
23:DB:1005:C:H2'	23:DB:1006:C:H6	1.72	0.54
23:DB:1083:U:H1'	23:DB:1086:A:N6	2.22	0.54
23:DB:1259:G:H2'	23:DB:1260:A:H8	1.73	0.54
23:DB:1637:A:H2'	23:DB:1638:C:H6	1.72	0.54
23:DB:184:C:H2'	23:DB:185:G:H8	1.71	0.54
23:DB:1904:G:O2'	23:DB:1905:C:H5'	2.07	0.54
23:DB:20:C:H2'	23:DB:21:A:H8	1.72	0.54
23:DB:554:U:H2'	23:DB:555:G:O4'	2.08	0.54
23:DB:593:U:H2'	23:DB:594:U:C6	2.42	0.54
23:DB:611:C:H2'	23:DB:612:G:O4'	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:696:G:O2'	23:DB:697:G:H5'	2.07	0.54
29:DG:29:ASN:HB2	29:DG:78:VAL:O	2.08	0.54
30:DH:2:GLN:HB2	30:DH:19:VAL:HA	1.89	0.54
30:DH:87:GLU:HB2	30:DH:89:LYS:HZ2	1.71	0.54
31:DJ:37:ARG:CZ	31:DJ:110:PRO:HG3	2.37	0.54
33:DL:58:TYR:HA	33:DL:62:PRO:HD3	1.89	0.54
34:DM:43:ALA:HB2	34:DM:69:PRO:HG3	1.89	0.54
41:DT:66:LYS:HD2	41:DT:67:VAL:N	2.22	0.54
44:DX:22:LEU:HD22	44:DX:25:GLN:OE1	2.07	0.54
46:DZ:47:LYS:HD2	46:DZ:52:ALA:HA	1.89	0.54
1:AA:235:C:H2'	1:AA:236:A:C8	2.43	0.54
1:AA:777:A:H2'	1:AA:778:G:H8	1.72	0.54
1:AA:971:G:OP1	1:AA:971:G:H3'	2.07	0.54
20:AB:186:VAL:HG22	20:AB:198:VAL:HG23	1.89	0.54
2:AC:149:LYS:HB3	2:AC:200:TRP:HB2	1.90	0.54
6:AG:67:ASN:ND2	6:AG:127:ALA:HA	2.20	0.54
8:AI:17:ARG:O	8:AI:64:ILE:HA	2.08	0.54
11:AL:87:LYS:NZ	11:AL:87:LYS:HA	2.21	0.54
12:AM:38:ILE:HG13	12:AM:55:LEU:HD21	1.89	0.54
12:AM:5:GLY:O	12:AM:6:ILE:HG12	2.07	0.54
13:AN:30:ILE:HG21	13:AN:44:VAL:HG21	1.89	0.54
13:AN:79:SER:O	13:AN:81:ILE:HD12	2.07	0.54
15:AP:31:ARG:HH11	15:AP:31:ARG:HG3	1.72	0.54
40:BS:38:TYR:CZ	47:B0:37:HIS:HA	2.42	0.54
49:B2:44:VAL:HG23	49:B2:45:SER:N	2.16	0.54
23:BB:2527:C:H1'	51:B4:1:MET:CB	2.38	0.54
23:BB:1121:C:H2'	23:BB:1122:G:O4'	2.07	0.54
23:BB:1480:C:H2'	23:BB:1481:U:C6	2.42	0.54
23:BB:1819:A:OP1	25:BC:153:LEU:HB2	2.08	0.54
23:BB:1821:A:H5'	25:BC:155:ARG:NH2	2.22	0.54
23:BB:2398:U:H2'	23:BB:2399:G:C8	2.42	0.54
23:BB:2840:C:H2'	23:BB:2841:C:H6	1.72	0.54
23:BB:288:U:O2'	23:BB:289:G:H5'	2.08	0.54
23:BB:633:A:C2'	23:BB:634:C:H5'	2.37	0.54
23:BB:802:A:H4'	55:BB:3474:HOH:O	2.08	0.54
25:BC:119:VAL:HG22	25:BC:130:PRO:HG2	1.89	0.54
26:BD:52:THR:HG22	26:BD:76:GLY:N	2.23	0.54
27:BE:141:MET:O	27:BE:143:LEU:HD22	2.08	0.54
31:BJ:15:TRP:HA	31:BJ:53:TYR:O	2.07	0.54
39:BR:5:PHE:HA	39:BR:40:MET:HG2	1.88	0.54
43:BW:47:GLY:O	43:BW:54:ARG:HA	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1131:G:C2'	1:CA:1132:C:H5'	2.36	0.54
1:CA:1524:C:H2'	1:CA:1525:G:C8	2.43	0.54
1:CA:22:G:H2'	1:CA:23:C:H6	1.73	0.54
20:CB:96:LEU:HB2	20:CB:99:MET:CG	2.38	0.54
12:CM:79:LEU:HD12	12:CM:80:MET:N	2.22	0.54
19:CT:2:ASN:O	19:CT:4:LYS:N	2.40	0.54
23:DB:136:G:H2'	23:DB:137:U:H6	1.69	0.54
23:DB:1430:G:H2'	23:DB:1431:A:C8	2.43	0.54
23:DB:1488:C:O2'	23:DB:1489:C:H5'	2.08	0.54
23:DB:1796:U:H2'	23:DB:1797:G:H8	1.73	0.54
23:DB:2271:G:O2'	23:DB:2272:U:H5'	2.07	0.54
26:DD:35:THR:HB	26:DD:48:ILE:CB	2.36	0.54
32:DK:115:ILE:HG23	32:DK:116:ILE:N	2.22	0.54
36:DO:18:LEU:HD13	43:DW:76:ARG:HH21	1.73	0.54
36:DO:90:VAL:HG12	36:DO:91:SER:N	2.23	0.54
26:DD:184:ARG:HD2	37:DP:4:ILE:CG2	2.38	0.54
39:DR:41:ILE:HG23	39:DR:43:ASN:HB2	1.88	0.54
43:DW:19:ARG:CZ	43:DW:19:ARG:HB2	2.37	0.54
43:DW:56:HIS:CD2	43:DW:57:THR:H	2.24	0.54
44:DX:1:MET:CB	44:DX:6:LEU:HG	2.37	0.54
45:DY:2:LYS:H	45:DY:37:ARG:CB	2.20	0.54
1:AA:1220:G:H2'	1:AA:1221:G:H8	1.72	0.54
1:AA:1380:U:O4	6:AG:2:ARG:HA	2.07	0.54
1:AA:230:G:O2'	1:AA:231:U:H5'	2.07	0.54
1:AA:936:C:H1'	1:AA:1382:C:H42	1.72	0.54
2:AC:46:LEU:HD11	2:AC:86:LEU:HD21	1.90	0.54
2:AC:78:LYS:CG	2:AC:81:GLU:HB2	2.34	0.54
4:AE:156:ARG:HD3	7:AH:42:GLU:O	2.07	0.54
8:AI:99:LYS:HD3	8:AI:100:ALA:N	2.21	0.54
8:AI:38:PHE:O	8:AI:44:ARG:HG2	2.08	0.54
8:AI:51:LEU:HD23	8:AI:56:MET:HE1	1.90	0.54
15:AP:8:ARG:HE	15:AP:15:PRO:HB3	1.72	0.54
15:AP:33:ILE:HD12	15:AP:33:ILE:N	2.21	0.54
50:B3:59:ALA:O	50:B3:60:CYS:HB2	2.07	0.54
23:BB:1024:G:C3'	23:BB:1025:G:H5''	2.34	0.54
23:BB:1166:G:H2'	23:BB:1167:C:H6	1.73	0.54
23:BB:1639:C:H2'	23:BB:1640:A:C5'	2.33	0.54
23:BB:272:A:H2'	23:BB:273:G:H8	1.72	0.54
23:BB:2843:G:O2'	23:BB:2844:G:H5'	2.08	0.54
23:BB:611:C:H2'	23:BB:612:G:O4'	2.07	0.54
23:BB:926:G:H2'	23:BB:927:A:C8	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:12:THR:OG1	26:BD:14:ILE:HG12	2.06	0.54
29:BG:172:GLU:O	29:BG:173:ALA:HB2	2.07	0.54
30:BH:4:ILE:C	30:BH:37:VAL:HG12	2.27	0.54
52:BI:23:VAL:HG23	52:BI:24:GLY:H	1.71	0.54
36:BO:16:ARG:HG3	36:BO:20:GLU:HG3	1.90	0.54
41:BT:32:LEU:C	41:BT:34:VAL:H	2.11	0.54
1:CA:1030:U:O2'	1:CA:1031:C:P	2.65	0.54
1:CA:1118:U:O2'	1:CA:1119:C:H5'	2.07	0.54
1:CA:128:G:H2'	1:CA:129:A:C8	2.43	0.54
1:CA:173:U:H5'	1:CA:197:A:O4'	2.07	0.54
1:CA:415:A:H3'	1:CA:416:G:H8	1.73	0.54
1:CA:499:A:H4'	1:CA:500:G:OP1	2.06	0.54
1:CA:979:C:H2'	1:CA:980:C:O4'	2.07	0.54
2:CC:133:MET:HE1	2:CC:165:GLU:HB3	1.89	0.54
3:CD:160:LEU:HD22	3:CD:160:LEU:O	2.07	0.54
5:CF:6:ILE:HD12	5:CF:7:VAL:N	2.22	0.54
8:CI:51:LEU:HB3	8:CI:56:MET:CG	2.34	0.54
9:CJ:65:TYR:HE1	13:CN:84:ARG:HA	1.72	0.54
9:CJ:65:TYR:OH	13:CN:84:ARG:HG3	2.07	0.54
11:CL:49:ARG:CG	11:CL:89:LEU:HD21	2.36	0.54
1:CA:1216:A:H5''	13:CN:4:SER:HB2	1.90	0.54
15:CP:40:ASN:ND2	15:CP:42:ILE:HG12	2.22	0.54
21:CU:4:LYS:HB3	21:CU:6:ARG:NH1	2.12	0.54
23:DB:1287:A:O2'	23:DB:1288:G:H5'	2.08	0.54
23:DB:1480:C:H2'	23:DB:1481:U:C6	2.43	0.54
23:DB:2070:A:H2'	23:DB:2071:A:C8	2.43	0.54
23:DB:2398:U:H2'	23:DB:2399:G:C8	2.42	0.54
23:DB:2556:C:H2'	23:DB:2557:G:O4'	2.08	0.54
23:DB:2839:G:H4'	35:DN:49:GLU:CG	2.30	0.54
23:DB:354:A:H2'	23:DB:355:U:H6	1.69	0.54
23:DB:512:G:H4'	23:DB:512:G:OP1	2.07	0.54
23:DB:580:U:H2'	23:DB:581:C:H6	1.70	0.54
23:DB:607:U:O4	23:DB:620:G:H5''	2.07	0.54
23:DB:70:G:H3'	23:DB:113:U:H4'	1.89	0.54
23:DB:851:C:H2'	23:DB:852:U:H6	1.73	0.54
26:DD:83:ARG:O	26:DD:84:LEU:HB2	2.07	0.54
27:DE:193:VAL:HA	27:DE:198:GLU:C	2.28	0.54
27:DE:75:SER:OG	27:DE:77:ILE:HG22	2.07	0.54
23:DB:2405:G:H4'	33:DL:70:LYS:HE3	1.90	0.54
33:DL:92:LEU:HA	33:DL:96:LYS:CD	2.37	0.54
41:DT:15:HIS:H	41:DT:32:LEU:HA	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:454:G:O2'	1:AA:455:G:H5'	2.07	0.54
20:AB:53:LEU:HD21	20:AB:212:TYR:CE2	2.42	0.54
2:AC:183:TYR:HA	2:AC:199:VAL:O	2.07	0.54
2:AC:96:VAL:HB	2:AC:97:PRO:HD2	1.88	0.54
3:AD:25:ARG:NH1	3:AD:30:LYS:HG2	2.23	0.54
13:AN:30:ILE:CD1	13:AN:30:ILE:H	2.18	0.54
13:AN:66:THR:HG23	13:AN:67:GLY:N	2.22	0.54
47:B0:42:ILE:HG12	47:B0:43:THR:N	2.22	0.54
23:BB:250:G:OP2	50:B3:7:ARG:HD3	2.08	0.54
22:BA:15:A:H1'	22:BA:109:A:C8	2.43	0.54
23:BB:1013:C:H2'	23:BB:1014:A:C8	2.42	0.54
23:BB:1097:U:C2'	23:BB:1098:A:H5'	2.36	0.54
23:BB:1559:U:H3'	23:BB:1560:G:H5'	1.88	0.54
23:BB:1667:G:OP1	32:BK:6:THR:HA	2.07	0.54
23:BB:1946:U:H2'	23:BB:1947:C:C6	2.43	0.54
23:BB:2218:G:H2'	23:BB:2219:U:C6	2.43	0.54
23:BB:2556:C:H2'	23:BB:2557:G:O4'	2.08	0.54
23:BB:2704:C:H2'	23:BB:2705:A:O4'	2.07	0.54
23:BB:36:G:H2'	23:BB:37:C:O4'	2.08	0.54
23:BB:233:A:N6	23:BB:428:A:H61	2.06	0.54
23:BB:480:A:H4'	42:BU:42:LYS:CG	2.32	0.54
23:BB:649:G:N2	50:B3:44:ARG:HH22	2.06	0.54
23:BB:705:A:H61	23:BB:726:G:H1'	1.73	0.54
23:BB:718:A:H3'	23:BB:719:C:C6	2.41	0.54
23:BB:877:A:C2	23:BB:900:A:N7	2.76	0.54
25:BC:124:LYS:O	25:BC:191:LEU:HD13	2.07	0.54
30:BH:73:ASN:HD22	30:BH:74:ALA:H	1.55	0.54
35:BN:107:ASN:HD21	40:BS:40:ASN:CG	2.11	0.54
35:BN:29:VAL:HB	35:BN:75:ILE:HG12	1.89	0.54
36:BO:41:ALA:HB3	36:BO:45:SER:O	2.07	0.54
38:BQ:88:GLU:HA	39:BR:53:PHE:CE1	2.42	0.54
41:BT:30:ILE:CG2	41:BT:31:VAL:N	2.70	0.54
42:BU:33:VAL:HG12	42:BU:35:VAL:CG2	2.38	0.54
1:CA:266:G:HO2'	1:CA:267:C:H3'	1.73	0.54
1:CA:337:G:H2'	1:CA:338:A:C8	2.43	0.54
1:CA:539:A:H2'	1:CA:540:G:H8	1.73	0.54
1:CA:777:A:H2'	1:CA:778:G:H8	1.72	0.54
1:CA:930:C:H2'	1:CA:931:C:C6	2.41	0.54
20:CB:86:CYS:HB3	20:CB:88:GLN:HG3	1.90	0.54
3:CD:151:GLN:HE21	3:CD:153:ARG:HD2	1.73	0.54
3:CD:2:ARG:HA	3:CD:2:ARG:HE	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:55:ARG:NH1	3:CD:55:ARG:HG3	2.20	0.54
6:CG:136:LYS:O	6:CG:140:VAL:HG23	2.07	0.54
6:CG:22:LEU:O	6:CG:26:VAL:HG13	2.06	0.54
1:CA:108:G:C6	19:CT:9:ARG:HG2	2.41	0.54
22:DA:102:G:H2'	22:DA:103:U:H6	1.72	0.54
23:DB:1168:G:O2'	23:DB:1169:A:H5'	2.07	0.54
23:DB:1487:U:H2'	23:DB:1488:C:H6	1.73	0.54
23:DB:1591:A:H2'	23:DB:1592:C:C6	2.42	0.54
23:DB:160:A:H2'	23:DB:161:A:C8	2.42	0.54
23:DB:431:U:O2'	23:DB:432:A:H5'	2.08	0.54
23:DB:535:G:O2'	23:DB:536:G:H5'	2.08	0.54
23:DB:95:A:O2'	44:DX:41:HIS:HA	2.08	0.54
23:DB:2025:C:P	26:DD:154:LYS:HE3	2.47	0.54
26:DD:60:VAL:HG23	26:DD:63:PRO:CG	2.33	0.54
52:DI:5:GLN:O	52:DI:6:ALA:HB3	2.07	0.54
34:DM:96:ILE:HG22	34:DM:97:GLN:NE2	2.22	0.54
39:DR:40:MET:CG	39:DR:54:VAL:HG13	2.36	0.54
1:AA:1281:C:H5'	1:AA:1282:C:H5	1.73	0.54
1:AA:1391:U:H2'	1:AA:1392:G:H8	1.72	0.54
1:AA:1452:C:H4'	1:AA:1453:G:C5'	2.37	0.54
1:AA:1527:U:O2'	1:AA:1528:U:H5'	2.07	0.54
2:AC:40:GLN:NE2	2:AC:44:LYS:HD2	2.22	0.54
1:AA:405:U:O4	3:AD:1:ALA:HA	2.07	0.54
11:AL:106:VAL:CG2	11:AL:116:TYR:HB3	2.38	0.54
15:AP:46:LYS:HD3	15:AP:46:LYS:H	1.73	0.54
19:AT:49:ALA:O	19:AT:52:GLU:HB3	2.08	0.54
48:B1:9:LYS:HG2	50:B3:34:LYS:NZ	2.23	0.54
23:BB:1210:G:H5''	23:BB:1211:C:O5'	2.07	0.54
23:BB:1430:G:H2'	23:BB:1431:A:C8	2.43	0.54
23:BB:1593:A:H2'	23:BB:1594:U:C6	2.42	0.54
23:BB:770:G:O2'	23:BB:771:G:H5'	2.07	0.54
25:BC:141:HIS:CG	25:BC:190:THR:HB	2.42	0.54
25:BC:67:LYS:CG	25:BC:68:ARG:H	2.14	0.54
26:BD:128:ARG:HD3	26:BD:142:VAL:O	2.08	0.54
26:BD:58:ASN:O	26:BD:59:ARG:HB2	2.07	0.54
27:BE:43:THR:O	27:BE:89:PRO:HA	2.07	0.54
32:BK:119:ALA:N	32:BK:120:PRO:HD3	2.23	0.54
33:BL:117:THR:HG21	33:BL:120:VAL:CB	2.37	0.54
34:BM:73:ILE:HD11	34:BM:92:TRP:HB2	1.89	0.54
35:BN:12:ARG:HD3	35:BN:20:MET:HG3	1.90	0.54
35:BN:83:LEU:HA	35:BN:86:ARG:HD3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BO:91:SER:HB2	36:BO:115:LEU:HG	1.89	0.54
37:BP:61:ARG:NH2	37:BP:75:THR:HB	2.22	0.54
23:BB:533:G:H5'	38:BQ:23:TYR:CD1	2.43	0.54
31:BJ:41:LYS:O	38:BQ:66:ALA:HB1	2.08	0.54
40:BS:86:MET:HB2	40:BS:96:ILE:CD1	2.38	0.54
41:BT:84:TYR:O	41:BT:85:VAL:HG22	2.08	0.54
24:BV:63:ILE:O	24:BV:65:VAL:HG22	2.07	0.54
1:CA:1296:C:H4'	1:CA:1302:C:C4	2.43	0.54
1:CA:1487:G:O2'	1:CA:1488:G:H5'	2.07	0.54
1:CA:766:A:H2	1:CA:1525:G:N3	2.06	0.54
30:BH:83:LYS:HG3	1:CA:360:G:OP1	2.07	0.54
1:CA:476:U:H2'	1:CA:477:C:H6	1.73	0.54
2:CC:51:VAL:HG11	2:CC:67:ILE:HD11	1.89	0.54
3:CD:196:GLU:O	3:CD:199:ILE:HG23	2.06	0.54
3:CD:58:GLN:NE2	3:CD:58:GLN:HA	2.16	0.54
4:CE:19:ARG:CG	4:CE:20:VAL:H	2.13	0.54
23:DB:2666:C:O4'	23:DB:2666:C:O2	2.24	0.54
23:DB:2692:G:O2'	23:DB:2693:G:H5'	2.07	0.54
23:DB:26:G:H2'	23:DB:27:G:C1'	2.37	0.54
23:DB:2704:C:H2'	23:DB:2705:A:O4'	2.07	0.54
23:DB:2849:U:H4'	23:DB:2850:A:C5'	2.37	0.54
23:DB:663:G:OP1	33:DL:27:LEU:HD13	2.08	0.54
27:DE:77:ILE:HG23	27:DE:78:TRP:HD1	1.73	0.54
30:DH:75:LEU:N	30:DH:75:LEU:HD23	2.22	0.54
31:DJ:23:LYS:C	31:DJ:25:LEU:HD13	2.28	0.54
35:DN:70:THR:HG22	35:DN:72:ASP:N	2.23	0.54
37:DP:78:PRO:O	37:DP:79:VAL:HG22	2.07	0.54
43:DW:4:LYS:HZ2	43:DW:4:LYS:H	1.54	0.54
1:AA:1244:G:H2'	1:AA:1245:C:H6	1.73	0.54
3:AD:18:LEU:HD22	3:AD:63:ILE:HG12	1.90	0.54
5:AF:18:VAL:N	5:AF:19:PRO:HD2	2.22	0.54
51:B4:24:ARG:HH11	51:B4:24:ARG:HG3	1.73	0.54
23:BB:1055:G:HO2'	23:BB:1085:A:H2	1.49	0.54
23:BB:1917:U:C2'	23:BB:1918:A:H5'	2.37	0.54
23:BB:329:G:H1	42:BU:16:LYS:HD3	1.73	0.54
23:BB:425:G:H2'	23:BB:426:C:H6	1.72	0.54
23:BB:581:C:H2'	23:BB:582:A:C8	2.43	0.54
23:BB:615:U:C2	27:BE:38:GLY:HA3	2.42	0.54
23:BB:741:U:H2'	23:BB:742:A:C8	2.42	0.54
23:BB:751:A:H5'	40:BS:90:LYS:CE	2.38	0.54
23:BB:936:A:H2'	23:BB:937:C:H6	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:124:LYS:CB	25:BC:125:PRO:CD	2.83	0.54
23:BB:1816:C:C5	25:BC:62:ARG:HD2	2.42	0.54
26:BD:131:ASP:O	26:BD:139:SER:HA	2.08	0.54
26:BD:175:LEU:HD13	26:BD:189:VAL:HG11	1.90	0.54
27:BE:119:ILE:N	27:BE:119:ILE:HD12	2.23	0.54
30:BH:114:GLU:OE2	30:BH:134:VAL:HG23	2.07	0.54
30:BH:37:VAL:O	30:BH:38:PRO:C	2.46	0.54
30:BH:83:LYS:HB2	30:BH:92:GLY:N	2.23	0.54
33:BL:82:LEU:O	33:BL:117:THR:HG22	2.06	0.54
33:BL:40:SER:C	33:BL:41:ARG:HE	2.10	0.54
33:BL:38:GLN:H	33:BL:41:ARG:HH11	1.56	0.54
39:BR:7:SER:C	39:BR:22:LEU:HD11	2.27	0.54
39:BR:47:VAL:HG22	39:BR:48:LYS:N	2.19	0.54
1:CA:235:C:H2'	1:CA:236:A:C8	2.43	0.54
1:CA:34:C:H2'	1:CA:35:G:C8	2.42	0.54
1:CA:719:C:H1'	17:CR:37:LYS:CB	2.32	0.54
20:CB:57:ASN:HD22	20:CB:223:GLY:HA2	1.73	0.54
3:CD:127:ARG:HD2	3:CD:127:ARG:H	1.73	0.54
8:CI:51:LEU:HD13	8:CI:56:MET:SD	2.48	0.54
13:CN:30:ILE:O	13:CN:44:VAL:HB	2.07	0.54
13:CN:26:LEU:HD21	13:CN:47:LEU:HG	1.90	0.54
1:CA:231:U:P	15:CP:31:ARG:HH12	2.31	0.54
23:DB:1387:A:H2'	23:DB:1388:G:C8	2.42	0.54
23:DB:1593:A:H2'	23:DB:1594:U:C6	2.42	0.54
23:DB:1723:G:N7	23:DB:1737:G:N2	2.56	0.54
23:DB:2213:U:O2	23:DB:2213:U:C2'	2.55	0.54
23:DB:2247:A:H2'	23:DB:2248:C:C6	2.43	0.54
23:DB:2257:U:C5'	43:DW:5:ALA:HB2	2.37	0.54
23:DB:2657:A:H2'	23:DB:2658:C:O4'	2.07	0.54
23:DB:533:G:H2'	23:DB:534:U:H6	1.73	0.54
26:DD:114:LYS:HB2	26:DD:114:LYS:HZ2	1.72	0.54
27:DE:102:ARG:HA	27:DE:105:LEU:HD23	1.90	0.54
27:DE:149:ILE:O	27:DE:149:ILE:HG13	2.08	0.54
27:DE:148:ILE:CA	27:DE:185:LYS:HB3	2.32	0.54
28:DF:141:ASP:HB3	28:DF:143:ASP:OD1	2.07	0.54
52:DI:121:ILE:CD1	52:DI:121:ILE:H	2.18	0.54
31:DJ:105:VAL:HG11	31:DJ:122:LEU:HD11	1.89	0.54
31:DJ:35:ARG:HH22	31:DJ:41:LYS:H	1.54	0.54
33:DL:47:ARG:O	33:DL:48:ARG:HD3	2.08	0.54
34:DM:91:TYR:N	34:DM:91:TYR:CD2	2.74	0.54
36:DO:26:LEU:O	36:DO:40:ILE:HD11	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:109:ILE:O	37:DP:111:GLU:N	2.40	0.54
38:DQ:39:ILE:HA	38:DQ:43:GLN:HE21	1.72	0.54
24:DV:53:LYS:NZ	24:DV:53:LYS:HB3	2.23	0.54
43:DW:19:ARG:HH21	43:DW:22:VAL:HA	1.73	0.54
23:DB:96:C:OP1	44:DX:41:HIS:HB2	2.08	0.54
46:DZ:14:ALA:HA	46:DZ:24:ILE:HA	1.89	0.54
46:DZ:49:ARG:C	46:DZ:51:VAL:N	2.61	0.54
1:AA:1244:G:H2'	1:AA:1245:C:C6	2.43	0.54
1:AA:1306:A:N6	1:AA:1331:G:H1'	2.23	0.54
1:AA:1477:U:H2'	1:AA:1478:U:C6	2.43	0.54
1:AA:255:G:H5'	16:AQ:17:GLU:O	2.08	0.54
1:AA:764:C:H2'	1:AA:765:G:H5'	1.88	0.54
12:AM:2:ARG:HG3	12:AM:8:ILE:HG23	1.89	0.54
13:AN:19:TYR:HD2	13:AN:23:ARG:HD3	1.73	0.54
48:B1:26:LYS:HD3	48:B1:29:LYS:HB2	1.88	0.54
23:BB:1723:G:N7	23:BB:1737:G:N2	2.56	0.54
23:BB:438:G:H2'	23:BB:439:A:C8	2.42	0.54
23:BB:709:U:H2'	23:BB:710:U:C6	2.43	0.54
23:BB:7:G:H2'	23:BB:8:C:O4'	2.08	0.54
25:BC:19:VAL:HG12	25:BC:20:ASN:N	2.15	0.54
23:BB:2811:G:H5'	26:BD:62:LYS:HD2	1.88	0.54
27:BE:180:LEU:HA	27:BE:186:VAL:HG21	1.88	0.54
27:BE:38:GLY:C	27:BE:40:ARG:H	2.09	0.54
23:BB:2761:A:H1'	29:BG:142:GLN:NE2	2.23	0.54
29:BG:84:LYS:NZ	29:BG:85:LYS:H	2.06	0.54
29:BG:94:ARG:CB	29:BG:94:ARG:HH11	2.20	0.54
52:BI:62:ALA:C	52:BI:64:ARG:H	2.10	0.54
32:BK:98:ARG:H	32:BK:98:ARG:HD2	1.71	0.54
34:BM:35:ALA:HB3	34:BM:97:GLN:C	2.28	0.54
35:BN:116:VAL:O	35:BN:117:ASP:HB2	2.08	0.54
39:BR:39:LEU:HD12	39:BR:60:LYS:N	2.23	0.54
23:BB:139:U:C4	41:BT:1:MET:HB2	2.42	0.54
43:BW:80:SER:OG	43:BW:84:GLU:HB2	2.08	0.54
46:BZ:30:HIS:O	46:BZ:31:ASP:HB3	2.07	0.54
1:CA:335:C:H2'	1:CA:336:A:H8	1.72	0.54
1:CA:845:A:H3'	1:CA:846:G:O4'	2.08	0.54
10:CK:83:VAL:HG21	10:CK:96:ILE:HG23	1.88	0.54
23:DB:1464:G:H2'	23:DB:1465:G:C8	2.43	0.54
23:DB:1515:A:H5'	23:DB:1557:C:H5'	1.88	0.54
23:DB:1439:A:C6	23:DB:1552:A:C5	2.95	0.54
23:DB:2020:A:H5'	47:D0:8:THR:HG23	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2250:G:C6	34:DM:81:ARG:HG2	2.42	0.54
23:DB:2386:A:H4'	43:DW:38:ARG:HB2	1.89	0.54
23:DB:2649:C:H2'	23:DB:2650:U:H6	1.73	0.54
26:DD:153:GLY:C	26:DD:155:VAL:H	2.11	0.54
26:DD:23:PRO:HA	26:DD:189:VAL:O	2.08	0.54
27:DE:97:ASN:HB2	27:DE:100:MET:HB2	1.89	0.54
52:DI:18:ASN:HB2	52:DI:38:CYS:SG	2.48	0.54
52:DI:89:SER:HA	52:DI:97:VAL:HG11	1.90	0.54
31:DJ:49:ASP:O	31:DJ:50:THR:CB	2.56	0.54
31:DJ:59:ALA:C	31:DJ:61:LYS:H	2.12	0.54
36:DO:9:ARG:C	36:DO:11:ALA:H	2.11	0.54
38:DQ:2:ARG:HG2	38:DQ:4:LYS:HE3	1.89	0.54
38:DQ:73:ILE:HG23	38:DQ:74:SER:N	2.22	0.54
41:DT:55:VAL:HG21	41:DT:85:VAL:HB	1.89	0.54
42:DU:28:LEU:HD13	42:DU:30:SER:N	2.07	0.54
42:DU:9:GLU:HB2	42:DU:71:ILE:CG1	2.38	0.54
24:DV:26:PHE:CE2	24:DV:44:HIS:HA	2.43	0.54
1:AA:1325:C:O2'	1:AA:1326:U:H5'	2.08	0.54
1:AA:337:G:H2'	1:AA:338:A:C8	2.43	0.54
1:AA:399:G:H2'	1:AA:400:C:C6	2.42	0.54
1:AA:472:U:H2'	1:AA:473:U:C6	2.43	0.54
3:AD:169:TRP:NE1	3:AD:170:LEU:HD23	2.22	0.54
4:AE:131:ASN:O	4:AE:135:VAL:HG22	2.08	0.54
4:AE:35:LEU:HD23	4:AE:36:THR:N	2.23	0.54
12:AM:2:ARG:HD2	12:AM:2:ARG:N	2.23	0.54
16:AQ:68:LYS:O	16:AQ:69:THR:CB	2.56	0.54
47:B0:36:LYS:HG3	47:B0:41:HIS:CA	2.38	0.54
48:B1:26:LYS:HD3	48:B1:29:LYS:HD3	1.90	0.54
23:BB:1061:U:H4'	23:BB:1070:A:O3'	2.08	0.54
23:BB:1083:U:H2'	23:BB:1085:A:OP2	2.08	0.54
23:BB:1171:G:H2'	23:BB:1172:C:C1'	2.38	0.54
23:BB:1287:A:O2'	23:BB:1288:G:H5'	2.08	0.54
23:BB:1400:U:H2'	23:BB:1401:G:C8	2.41	0.54
23:BB:1413:A:H2'	23:BB:1414:C:C6	2.43	0.54
23:BB:1716:U:H2'	23:BB:1717:A:C8	2.43	0.54
23:BB:1745:A:H2'	23:BB:1746:A:H8	1.73	0.54
23:BB:1797:G:OP1	25:BC:251:THR:HG23	2.08	0.54
23:BB:1934:C:O2'	23:BB:1935:G:H5'	2.08	0.54
23:BB:2489:U:O2'	23:BB:2490:G:H5'	2.08	0.54
23:BB:322:A:H5'	23:BB:340:A:C1'	2.38	0.54
23:BB:433:C:O2'	23:BB:434:U:H5'	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:34:VAL:HA	26:BD:89:GLU:CB	2.32	0.54
28:BF:59:ILE:HD13	28:BF:59:ILE:O	2.08	0.54
52:BI:14:ALA:HB1	52:BI:50:LYS:HA	1.89	0.54
23:BB:2780:G:N1	31:BJ:120:ARG:HG3	2.22	0.54
31:BJ:17:VAL:CG1	31:BJ:57:LEU:HD21	2.38	0.54
32:BK:61:VAL:HG23	32:BK:84:CYS:O	2.08	0.54
34:BM:59:ARG:N	34:BM:59:ARG:HD2	2.22	0.54
35:BN:10:LEU:O	35:BN:12:ARG:N	2.40	0.54
37:BP:99:LEU:HD21	37:BP:109:ILE:HG23	1.90	0.54
37:BP:23:ASP:HA	37:BP:48:ALA:O	2.08	0.54
38:BQ:52:ARG:NH2	38:BQ:55:GLN:HG2	2.23	0.54
44:BX:44:LYS:O	44:BX:44:LYS:HD2	2.08	0.54
45:BY:18:LYS:HD2	45:BY:53:MET:HE1	1.90	0.54
1:CA:1135:U:H4'	1:CA:1136:C:OP1	2.05	0.54
1:CA:472:U:H2'	1:CA:473:U:C6	2.43	0.54
4:CE:73:VAL:HG23	4:CE:146:MET:HE1	1.90	0.54
6:CG:110:ARG:HE	6:CG:122:GLU:HB2	1.72	0.54
14:CO:80:LEU:HD21	14:CO:84:LEU:HD22	1.90	0.54
19:CT:27:MET:HG2	19:CT:31:ILE:HD11	1.90	0.54
51:D4:7:VAL:O	51:D4:8:LYS:HB2	2.08	0.54
23:DB:1000:A:H2'	23:DB:1001:A:C8	2.43	0.54
23:DB:1219:U:H2'	23:DB:1220:G:H8	1.73	0.54
23:DB:1671:U:H2'	23:DB:1673:G:OP2	2.08	0.54
23:DB:2405:G:H1'	23:DB:2412:A:N6	2.23	0.54
23:DB:2557:G:H2'	23:DB:2558:C:H6	1.70	0.54
23:DB:2741:A:H2'	23:DB:2742:G:O4'	2.08	0.54
23:DB:2840:C:H2'	23:DB:2841:C:H6	1.72	0.54
23:DB:581:C:H2'	23:DB:582:A:C8	2.43	0.54
23:DB:770:G:O2'	23:DB:771:G:H5'	2.08	0.54
23:DB:1773:A:H62	25:DC:206:LYS:HE2	1.73	0.54
52:DI:10:LEU:O	52:DI:10:LEU:HD12	2.08	0.54
23:DB:1132:U:H5''	31:DJ:85:LYS:HD3	1.89	0.54
32:DK:24:VAL:HA	32:DK:39:ILE:CD1	2.38	0.54
32:DK:6:THR:HG22	32:DK:7:MET:N	2.23	0.54
34:DM:50:ARG:HA	34:DM:53:MET:HE2	1.90	0.54
36:DO:82:ALA:O	36:DO:83:LEU:HD12	2.08	0.54
37:DP:55:HIS:O	37:DP:57:ALA:N	2.38	0.54
23:DB:2387:U:O2'	43:DW:37:VAL:HG11	2.08	0.54
45:DY:1:ALA:O	45:DY:43:ILE:HG13	2.08	0.54
1:AA:1071:C:H2'	1:AA:1072:G:C8	2.40	0.54
1:AA:1461:G:H2'	1:AA:1462:C:H6	1.69	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:22:G:H2'	1:AA:23:C:H6	1.73	0.54
1:AA:471:U:H2'	1:AA:472:U:H6	1.72	0.54
1:AA:591:U:H2'	1:AA:592:G:H8	1.71	0.54
1:AA:736:C:H2'	1:AA:737:C:C6	2.43	0.54
1:AA:787:A:O2'	1:AA:788:U:H5'	2.08	0.54
20:AB:95:TRP:CE3	20:AB:171:ALA:HA	2.43	0.54
2:AC:100:ILE:C	2:AC:100:ILE:HD13	2.29	0.54
6:AG:14:ASP:OD1	6:AG:15:PRO:HD2	2.08	0.54
7:AH:29:SER:CB	7:AH:32:LYS:HZ2	2.21	0.54
16:AQ:74:LEU:HD22	16:AQ:75:VAL:N	2.22	0.54
18:AS:62:THR:N	18:AS:65:MET:HE3	2.23	0.54
1:AA:177:G:H5''	19:AT:59:ARG:HH21	1.73	0.54
22:BA:86:G:H2'	22:BA:87:U:O4'	2.08	0.54
23:BB:1149:G:H2'	23:BB:1150:C:C6	2.43	0.54
23:BB:1432:G:O2'	23:BB:1433:A:H5'	2.07	0.54
23:BB:1441:G:H2'	23:BB:1442:U:H6	1.73	0.54
23:BB:1464:G:H2'	23:BB:1465:G:C8	2.43	0.54
23:BB:1751:U:H2'	23:BB:1752:C:C6	2.43	0.54
23:BB:1682:G:C2	23:BB:1757:A:O4'	2.61	0.54
23:BB:2272:U:H3'	23:BB:2272:U:H6	1.73	0.54
23:BB:331:C:O2'	23:BB:332:A:H5'	2.08	0.54
23:BB:854:C:H2'	23:BB:855:G:C8	2.41	0.54
23:BB:855:G:H21	43:BW:23:LYS:HB3	1.70	0.54
23:BB:871:U:H2'	23:BB:872:U:C6	2.43	0.54
23:BB:956:G:H5''	34:BM:76:LYS:HE2	1.89	0.54
23:BB:992:C:O2'	23:BB:993:G:H5'	2.08	0.54
27:BE:198:GLU:C	27:BE:200:LEU:H	2.12	0.54
28:BF:65:LEU:C	28:BF:67:THR:H	2.11	0.54
52:BI:124:MET:O	52:BI:128:ILE:HG12	2.08	0.54
33:BL:119:PRO:O	33:BL:120:VAL:HB	2.06	0.54
34:BM:96:ILE:HD13	34:BM:97:GLN:CB	2.38	0.54
37:BP:32:VAL:HG13	37:BP:40:GLN:NE2	2.23	0.54
41:BT:61:LEU:HD12	41:BT:62:VAL:N	2.22	0.54
43:BW:43:LYS:O	43:BW:45:HIS:N	2.41	0.54
46:BZ:49:ARG:HD3	46:BZ:50:ASP:N	2.22	0.54
1:CA:1314:C:H2'	1:CA:1315:U:C6	2.43	0.54
1:CA:366:A:O2'	1:CA:367:U:OP1	2.24	0.54
1:CA:44:A:O2'	1:CA:45:G:H5'	2.08	0.54
1:CA:600:A:H2'	1:CA:601:G:H8	1.73	0.54
20:CB:46:VAL:CG1	20:CB:47:PRO:HD3	2.22	0.54
20:CB:78:ALA:CB	20:CB:213:LEU:HD13	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:115:GLN:NE2	3:CD:153:ARG:HH12	2.05	0.54
4:CE:32:PHE:CE2	4:CE:55:VAL:HG22	2.43	0.54
5:CF:18:VAL:HG21	5:CF:58:HIS:CE1	2.42	0.54
8:CI:23:GLY:H	8:CI:60:LEU:HA	1.73	0.54
12:CM:53:ASP:HA	12:CM:56:ARG:NH2	2.22	0.54
19:CT:79:THR:HA	19:CT:82:ILE:HG12	1.90	0.54
10:CK:121:ARG:HE	21:CU:34:ARG:CG	2.21	0.54
48:D1:22:THR:HB	50:D3:34:LYS:HE3	1.90	0.54
33:DL:59:ARG:HA	50:D3:11:LYS:NZ	2.22	0.54
23:DB:1745:A:H2'	23:DB:1746:A:H8	1.73	0.54
23:DB:2323:G:C2'	23:DB:2324:U:H5'	2.37	0.54
23:DB:345:A:H1'	23:DB:346:A:C2	2.41	0.54
25:DC:182:LYS:HE2	25:DC:264:LYS:NZ	2.23	0.54
23:DB:2024:G:H5''	26:DD:154:LYS:HZ3	1.73	0.54
52:DI:41:PHE:CE2	52:DI:45:THR:HG21	2.43	0.54
31:DJ:93:ILE:CG2	31:DJ:97:PRO:HA	2.37	0.54
35:DN:123:GLU:HG2	35:DN:124:ALA:H	1.72	0.54
35:DN:8:ARG:HD3	35:DN:43:GLU:OE2	2.07	0.54
36:DO:35:ILE:HG21	36:DO:71:ALA:HA	1.90	0.54
39:DR:80:ARG:HD2	39:DR:85:LYS:HA	1.90	0.54
42:DU:9:GLU:OE1	42:DU:71:ILE:HG13	2.08	0.54
43:DW:67:LYS:HG2	43:DW:71:LYS:CA	2.38	0.54
1:AA:916:U:H2'	1:AA:917:G:H8	1.72	0.53
7:AH:76:ARG:HG2	7:AH:77:VAL:N	2.23	0.53
47:B0:35:GLU:CD	47:B0:43:THR:HG22	2.28	0.53
47:B0:6:LYS:HZ2	47:B0:6:LYS:HA	1.72	0.53
23:BB:1488:C:O2'	23:BB:1489:C:H5'	2.08	0.53
23:BB:2199:A:H5'	23:BB:2200:C:OP2	2.07	0.53
23:BB:2292:U:H2'	23:BB:2293:G:C8	2.44	0.53
23:BB:636:G:O6	33:BL:76:GLU:HG2	2.08	0.53
23:BB:2750:A:H5''	29:BG:3:VAL:HG13	1.90	0.53
37:BP:112:ARG:HA	37:BP:112:ARG:HH11	1.73	0.53
37:BP:8:GLU:HA	37:BP:11:GLN:CG	2.37	0.53
41:BT:68:LYS:O	41:BT:68:LYS:HD3	2.08	0.53
41:BT:69:ARG:H	41:BT:75:GLY:HA3	1.72	0.53
1:CA:1245:C:H2'	1:CA:1246:A:H8	1.73	0.53
1:CA:930:C:H2'	1:CA:931:C:H6	1.73	0.53
20:CB:40:ILE:CG2	20:CB:200:PRO:HB2	2.38	0.53
20:CB:57:ASN:HB2	20:CB:219:THR:O	2.08	0.53
2:CC:113:LYS:HA	2:CC:184:ASN:ND2	2.22	0.53
3:CD:144:ILE:HD12	3:CD:154:VAL:HG11	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:51:ILE:HD13	5:CF:86:ARG:HG3	1.90	0.53
19:CT:27:MET:O	19:CT:31:ILE:HG13	2.07	0.53
47:D0:15:ARG:HG3	47:D0:15:ARG:HH11	1.73	0.53
48:D1:24:LYS:HB2	48:D1:24:LYS:HZ2	1.73	0.53
48:D1:35:LEU:HA	48:D1:48:TYR:O	2.08	0.53
22:DA:13:G:H4'	22:DA:15:A:H2'	1.89	0.53
23:DB:1413:A:H2'	23:DB:1414:C:C6	2.42	0.53
23:DB:1549:A:H2'	23:DB:1550:C:C6	2.43	0.53
23:DB:2216:G:H2'	23:DB:2217:G:C8	2.43	0.53
23:DB:2743:U:C2'	23:DB:2744:G:H5''	2.38	0.53
23:DB:873:C:H2'	23:DB:874:G:C8	2.43	0.53
26:DD:5:VAL:HG22	26:DD:51:THR:O	2.08	0.53
30:DH:121:VAL:O	30:DH:122:LEU:HB2	2.08	0.53
34:DM:4:PRO:HD3	34:DM:47:GLU:CG	2.39	0.53
34:DM:88:ASN:ND2	34:DM:88:ASN:N	2.56	0.53
37:DP:29:VAL:CG2	37:DP:47:ILE:HD11	2.37	0.53
39:DR:68:ARG:HB2	39:DR:97:LYS:HG3	1.90	0.53
43:DW:9:THR:OG1	43:DW:10:ARG:N	2.40	0.53
46:DZ:3:LYS:HB2	46:DZ:7:PRO:HA	1.90	0.53
3:AD:1:ALA:O	3:AD:67:LEU:HD21	2.08	0.53
4:AE:158:LYS:HE2	4:AE:158:LYS:H	1.73	0.53
6:AG:3:ARG:CZ	6:AG:3:ARG:HB3	2.37	0.53
10:AK:61:ALA:O	10:AK:64:VAL:HG12	2.08	0.53
10:AK:92:ARG:NH1	10:AK:92:ARG:HB3	2.20	0.53
12:AM:44:ILE:HD12	12:AM:45:SER:N	2.20	0.53
14:AO:60:SER:HA	14:AO:63:ARG:NH1	2.22	0.53
17:AR:33:THR:HG23	17:AR:35:SER:H	1.72	0.53
21:AU:29:ALA:HB1	21:AU:32:ARG:HH21	1.72	0.53
50:B3:11:LYS:HG2	50:B3:12:ARG:HG3	1.90	0.53
23:BB:1013:C:H2'	23:BB:1014:A:H8	1.73	0.53
23:BB:1139:G:O2'	23:BB:1140:C:H5'	2.09	0.53
23:BB:1259:G:H2'	23:BB:1260:A:C8	2.42	0.53
23:BB:139:U:C5	41:BT:1:MET:HB2	2.44	0.53
23:BB:2204:G:H5'	25:BC:149:LYS:CE	2.25	0.53
23:BB:418:C:H2'	23:BB:419:U:H6	1.73	0.53
25:BC:28:PRO:HD2	25:BC:79:ARG:CZ	2.37	0.53
27:BE:91:ASP:O	27:BE:92:HIS:HB2	2.06	0.53
28:BF:3:LEU:O	28:BF:3:LEU:HD23	2.08	0.53
32:BK:36:GLY:HA2	32:BK:62:VAL:O	2.09	0.53
34:BM:67:VAL:CG2	34:BM:95:LEU:HD21	2.39	0.53
35:BN:9:GLN:HE21	35:BN:17:ARG:NH1	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:52:ILE:HG21	35:BN:94:TYR:CD2	2.43	0.53
35:BN:76:VAL:O	35:BN:79:LEU:HB2	2.08	0.53
40:BS:71:VAL:CG2	40:BS:107:VAL:HG12	2.30	0.53
40:BS:29:VAL:CG1	40:BS:69:LEU:HB2	2.38	0.53
1:CA:1001:C:H2'	1:CA:1002:G:H8	1.74	0.53
1:CA:1347:G:N2	1:CA:1373:G:H2'	2.24	0.53
1:CA:1468:A:O2'	1:CA:1469:C:H5'	2.08	0.53
1:CA:961:U:O4'	1:CA:961:U:O2	2.26	0.53
3:CD:22:SER:H	3:CD:109:THR:HG22	1.72	0.53
3:CD:36:ALA:C	3:CD:38:GLY:H	2.11	0.53
3:CD:95:GLY:HA3	3:CD:135:GLN:NE2	2.24	0.53
8:CI:62:LEU:N	8:CI:62:LEU:HD13	2.23	0.53
15:CP:67:ILE:HG12	15:CP:72:ALA:HB2	1.91	0.53
49:D2:13:ASN:C	49:D2:15:SER:H	2.10	0.53
33:DL:58:TYR:HB3	50:D3:13:PHE:CE1	2.42	0.53
22:DA:32:U:H4'	22:DA:52:A:N6	2.23	0.53
23:DB:2186:G:O2'	23:DB:2187:U:H5'	2.07	0.53
23:DB:251:A:H2'	23:DB:252:G:O4'	2.08	0.53
23:DB:27:G:N2	23:DB:512:G:H2'	2.16	0.53
23:DB:65:U:H2'	23:DB:66:C:C6	2.42	0.53
26:DD:17:GLU:OE2	32:DK:72:PRO:HB2	2.08	0.53
28:DF:9:ASP:O	28:DF:10:GLU:HB2	2.08	0.53
28:DF:137:PHE:O	28:DF:139:GLU:N	2.41	0.53
23:DB:2311:A:N3	28:DF:39:VAL:HB	2.22	0.53
30:DH:122:LEU:HD13	30:DH:146:VAL:HG13	1.90	0.53
31:DJ:100:VAL:HG22	31:DJ:101:ILE:H	1.72	0.53
37:DP:90:ALA:HB3	37:DP:111:GLU:CB	2.38	0.53
39:DR:69:GLY:CA	39:DR:96:VAL:HG13	2.39	0.53
46:DZ:24:ILE:CD1	46:DZ:24:ILE:H	2.14	0.53
1:AA:1130:A:H5'	8:AI:19:PHE:CE2	2.43	0.53
1:AA:1237:C:H4'	1:AA:1334:G:N2	2.24	0.53
1:AA:203:G:N2	1:AA:205:A:N6	2.57	0.53
1:AA:474:G:H2'	1:AA:475:C:H6	1.73	0.53
1:AA:730:G:O2'	1:AA:766:A:H5'	2.08	0.53
20:AB:79:VAL:HG11	20:AB:92:ASN:HB3	1.89	0.53
2:AC:110:LEU:O	2:AC:203:LYS:HE2	2.08	0.53
6:AG:144:ALA:C	6:AG:146:ALA:H	2.11	0.53
8:AI:71:ILE:CD1	8:AI:71:ILE:H	2.20	0.53
8:AI:91:GLU:CD	8:AI:94:ARG:HH21	2.11	0.53
9:AJ:67:ILE:HG12	13:AN:94:GLY:O	2.08	0.53
15:AP:40:ASN:HD21	15:AP:42:ILE:HG13	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1231:U:H2'	23:BB:1232:G:H8	1.73	0.53
23:BB:1204:A:N1	23:BB:1241:A:N1	2.55	0.53
23:BB:152:A:H2'	23:BB:153:U:C6	2.43	0.53
23:BB:155:A:H2'	23:BB:156:A:H8	1.68	0.53
23:BB:1629:U:O2	23:BB:2698:U:H5''	2.08	0.53
23:BB:2053:G:C2'	23:BB:2054:A:H5'	2.37	0.53
23:BB:2774:C:H2'	23:BB:2775:G:O4'	2.09	0.53
23:BB:2848:G:H1'	23:BB:2868:A:H61	1.72	0.53
23:BB:696:G:O2'	23:BB:697:G:H5'	2.07	0.53
23:BB:711:G:O2'	23:BB:712:G:H5'	2.09	0.53
23:BB:877:A:H2'	23:BB:899:A:N1	2.24	0.53
26:BD:59:ARG:HH21	26:BD:60:VAL:CG1	2.22	0.53
26:BD:59:ARG:HH21	26:BD:60:VAL:HG12	1.73	0.53
26:BD:62:LYS:N	26:BD:63:PRO:HD2	2.23	0.53
27:BE:144:GLU:C	27:BE:146:VAL:H	2.11	0.53
29:BG:17:LYS:HE3	29:BG:24:THR:CG2	2.37	0.53
30:BH:1:MET:HB2	30:BH:20:ASN:OD1	2.09	0.53
52:BI:122:GLU:CD	52:BI:122:GLU:H	2.12	0.53
52:BI:23:VAL:HG23	52:BI:24:GLY:N	2.24	0.53
52:BI:37:PHE:HB2	52:BI:66:PHE:CZ	2.44	0.53
31:BJ:72:LYS:HB3	31:BJ:74:TYR:CE2	2.43	0.53
24:BV:84:PRO:HG3	34:BM:127:LYS:CE	2.37	0.53
34:BM:3:GLN:O	34:BM:4:PRO:C	2.46	0.53
34:BM:73:ILE:HG22	34:BM:74:THR:N	2.24	0.53
37:BP:69:VAL:O	37:BP:70:GLU:HB2	2.09	0.53
39:BR:41:ILE:C	39:BR:43:ASN:N	2.60	0.53
42:BU:90:LYS:NZ	42:BU:91:LYS:HE2	2.23	0.53
46:BZ:25:ARG:HG3	46:BZ:26:SER:N	2.24	0.53
1:CA:1252:A:H2'	1:CA:1253:G:O4'	2.08	0.53
1:CA:317:U:H2'	1:CA:318:G:H8	1.73	0.53
1:CA:672:U:H2'	1:CA:673:A:C8	2.43	0.53
1:CA:677:U:H2'	1:CA:678:U:C6	2.44	0.53
1:CA:860:A:H2'	1:CA:861:G:O4'	2.09	0.53
2:CC:116:ALA:HB1	2:CC:186:SER:OG	2.07	0.53
1:CA:545:C:H5''	3:CD:68:GLU:HG2	1.90	0.53
5:CF:10:VAL:HG12	5:CF:11:HIS:N	2.23	0.53
6:CG:14:ASP:HB3	6:CG:18:GLY:H	1.73	0.53
9:CJ:82:LYS:HD2	9:CJ:82:LYS:N	2.23	0.53
10:CK:70:ALA:C	10:CK:72:ALA:H	2.11	0.53
18:CS:15:LEU:HD23	18:CS:15:LEU:O	2.08	0.53
23:DB:1100:C:H2'	23:DB:1101:U:C6	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1172:C:H2'	23:DB:1173:U:O4'	2.07	0.53
23:DB:1682:G:C2	23:DB:1757:A:O4'	2.61	0.53
23:DB:1716:U:H2'	23:DB:1717:A:C8	2.43	0.53
23:DB:2077:A:O2'	23:DB:2078:C:H5'	2.08	0.53
23:DB:6:A:H4'	31:DJ:133:ALA:O	2.08	0.53
23:DB:905:A:O2'	23:DB:906:U:H5'	2.09	0.53
25:DC:22:GLU:HB2	25:DC:202:ARG:CG	2.38	0.53
25:DC:245:THR:O	25:DC:247:TRP:N	2.40	0.53
28:DF:31:GLU:O	28:DF:95:MET:HE1	2.08	0.53
29:DG:171:LYS:HD3	29:DG:172:GLU:C	2.29	0.53
29:DG:25:ILE:HD11	29:DG:75:VAL:HG22	1.89	0.53
23:DB:1080:A:O2'	52:DI:126:ARG:HD2	2.08	0.53
32:DK:43:ILE:HD11	32:DK:58:LEU:HD21	1.90	0.53
33:DL:96:LYS:HE3	33:DL:96:LYS:CA	2.38	0.53
33:DL:33:ARG:HG2	39:DR:85:LYS:HZ3	1.73	0.53
40:DS:74:ILE:CG2	40:DS:105:VAL:HG23	2.38	0.53
1:AA:128:G:H2'	1:AA:129:A:C8	2.43	0.53
1:AA:1314:C:H2'	1:AA:1315:U:C6	2.44	0.53
1:AA:817:C:H1'	1:AA:819:A:C5'	2.38	0.53
1:AA:982:U:H4'	1:AA:983:A:O4'	2.08	0.53
7:AH:42:GLU:HG3	7:AH:100:ILE:HD13	1.90	0.53
8:AI:30:ASN:HD22	8:AI:30:ASN:N	2.06	0.53
12:AM:52:ILE:HD12	12:AM:55:LEU:HD12	1.90	0.53
14:AO:11:VAL:HA	14:AO:26:VAL:HG13	1.90	0.53
14:AO:24:THR:HB	14:AO:69:LEU:HD21	1.90	0.53
16:AQ:23:ALA:C	16:AQ:24:ILE:HD12	2.29	0.53
18:AS:12:LEU:HG	18:AS:15:LEU:HD23	1.91	0.53
22:BA:35:C:H2'	22:BA:36:C:O4'	2.08	0.53
23:BB:1159:U:H2'	23:BB:1160:G:C8	2.43	0.53
23:BB:1387:A:H2'	23:BB:1388:G:C8	2.41	0.53
23:BB:1439:A:N6	23:BB:1440:U:O2	2.35	0.53
23:BB:1487:U:H2'	23:BB:1488:C:H6	1.73	0.53
23:BB:1571:A:H2'	23:BB:1572:A:H8	1.73	0.53
23:BB:1733:G:H2'	23:BB:1734:G:H8	1.74	0.53
23:BB:1842:G:H2'	23:BB:1843:C:C6	2.44	0.53
23:BB:2305:U:H1'	28:BF:132:ARG:HA	1.90	0.53
23:BB:2405:G:H1'	23:BB:2412:A:N6	2.23	0.53
23:BB:2772:C:H5''	26:BD:171:THR:HG21	1.91	0.53
23:BB:852:U:H2'	23:BB:853:C:H6	1.72	0.53
23:BB:987:C:H2'	23:BB:988:A:O4'	2.08	0.53
23:BB:992:C:H2'	23:BB:993:G:H8	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:155:ARG:NE	25:BC:157:ALA:HB3	2.16	0.53
23:BB:2636:C:H4'	26:BD:80:TRP:CE3	2.43	0.53
29:BG:78:VAL:HG23	29:BG:79:THR:HG23	1.91	0.53
31:BJ:43:GLU:O	31:BJ:45:THR:N	2.42	0.53
31:BJ:84:ILE:HG22	31:BJ:85:LYS:H	1.73	0.53
33:BL:64:PHE:H	50:B3:24:LYS:CD	2.14	0.53
24:BV:57:TYR:CD2	34:BM:133:LYS:HE2	2.43	0.53
34:BM:95:LEU:HD22	34:BM:98:PRO:CG	2.36	0.53
36:BO:8:ILE:C	36:BO:10:ARG:H	2.12	0.53
37:BP:108:ARG:CD	37:BP:108:ARG:H	2.21	0.53
38:BQ:27:ARG:HA	38:BQ:33:VAL:HB	1.90	0.53
39:BR:62:GLU:HG3	39:BR:103:ALA:HA	1.89	0.53
1:CA:1412:C:H2'	1:CA:1413:A:H8	1.72	0.53
1:CA:787:A:O2'	1:CA:788:U:H5'	2.08	0.53
20:CB:62:ARG:H	20:CB:62:ARG:CD	2.22	0.53
20:CB:86:CYS:HB3	20:CB:88:GLN:NE2	2.22	0.53
4:CE:104:ILE:HG13	4:CE:114:LEU:HB2	1.89	0.53
8:CI:24:ASN:OD1	8:CI:26:LYS:HG2	2.09	0.53
12:CM:95:PRO:HD3	12:CM:108:ARG:HG2	1.88	0.53
18:CS:28:LYS:HB3	18:CS:28:LYS:NZ	2.23	0.53
51:D4:23:ILE:HD12	51:D4:24:ARG:HD2	1.90	0.53
51:D4:2:LYS:HG2	51:D4:38:GLY:HA3	1.90	0.53
22:DA:21:G:H2'	22:DA:22:U:O4'	2.08	0.53
22:DA:53:A:O2'	22:DA:54:G:H5'	2.07	0.53
23:DB:1051:G:H2'	23:DB:1052:C:H6	1.73	0.53
23:DB:1198:U:C2	23:DB:1199:U:C5	2.97	0.53
23:DB:130:C:O2'	23:DB:131:A:H5'	2.08	0.53
23:DB:1751:U:H2'	23:DB:1752:C:C6	2.43	0.53
23:DB:2247:A:H2'	23:DB:2248:C:H6	1.71	0.53
23:DB:247:G:H4'	23:DB:386:G:C4	2.43	0.53
23:DB:2874:C:H2'	23:DB:2875:C:C6	2.42	0.53
23:DB:545:U:H3'	23:DB:546:U:H5''	1.91	0.53
26:DD:79:LEU:CG	26:DD:80:TRP:H	2.20	0.53
27:DE:191:ASP:HA	27:DE:194:LYS:CE	2.39	0.53
28:DF:168:LEU:HD23	28:DF:168:LEU:O	2.09	0.53
29:DG:39:ALA:HA	29:DG:54:ARG:HB2	1.90	0.53
30:DH:90:LEU:HD22	30:DH:122:LEU:O	2.07	0.53
30:DH:95:GLY:O	30:DH:99:ILE:HG12	2.08	0.53
52:DI:102:ARG:HG3	52:DI:141:ASP:CB	2.38	0.53
31:DJ:23:LYS:HE3	31:DJ:63:ALA:CB	2.38	0.53
34:DM:86:LYS:HG2	34:DM:87:GLY:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:32:VAL:HA	37:DP:42:PHE:CB	2.34	0.53
41:DT:6:ARG:HG2	41:DT:9:LYS:O	2.09	0.53
43:DW:30:VAL:HG12	43:DW:31:LEU:N	2.23	0.53
45:DY:1:ALA:N	45:DY:43:ILE:HB	2.23	0.53
1:AA:610:U:O2	1:AA:610:U:O4'	2.25	0.53
1:AA:845:A:H3'	1:AA:846:G:O4'	2.08	0.53
20:AB:165:ALA:CB	20:AB:186:VAL:HG12	2.38	0.53
23:BB:254:G:O6	50:B3:3:ILE:HG23	2.09	0.53
23:BB:1220:G:H2'	23:BB:1221:C:H6	1.72	0.53
23:BB:2304:G:H22	23:BB:2312:U:H3	1.56	0.53
23:BB:2794:C:H2'	23:BB:2795:C:C6	2.44	0.53
23:BB:520:G:H2'	23:BB:521:U:C6	2.43	0.53
23:BB:716:A:H3'	23:BB:717:C:H5''	1.90	0.53
27:BE:44:ARG:H	27:BE:44:ARG:CD	2.21	0.53
29:BG:26:LYS:HG2	29:BG:31:GLU:HG2	1.90	0.53
29:BG:90:GLY:HA3	29:BG:93:TYR:CD2	2.44	0.53
34:BM:67:VAL:HG23	34:BM:95:LEU:HD21	1.89	0.53
40:BS:69:LEU:N	40:BS:69:LEU:HD22	2.24	0.53
24:BV:82:TYR:CD2	24:BV:82:TYR:N	2.76	0.53
1:CA:633:G:H2'	1:CA:634:C:C6	2.44	0.53
1:CA:920:U:H2'	1:CA:921:U:C6	2.43	0.53
5:CF:51:ILE:CD1	5:CF:86:ARG:HG3	2.39	0.53
15:CP:44:SER:C	15:CP:46:LYS:H	2.12	0.53
50:D3:24:LYS:HZ2	50:D3:24:LYS:HB3	1.74	0.53
23:DB:1545:A:H2'	23:DB:1546:G:O4'	2.09	0.53
23:DB:1842:G:H2'	23:DB:1843:C:C6	2.43	0.53
23:DB:2723:C:H5'	35:DN:4:ARG:NH2	2.24	0.53
25:DC:28:PRO:O	25:DC:29:PHE:CG	2.61	0.53
23:DB:1813:G:H1'	25:DC:45:ASN:HB3	1.90	0.53
25:DC:53:ILE:CG1	25:DC:218:THR:HA	2.37	0.53
26:DD:124:ARG:NH1	26:DD:160:LYS:HE2	2.24	0.53
26:DD:173:GLN:HG2	26:DD:208:LYS:HB3	1.90	0.53
29:DG:169:ARG:CZ	29:DG:169:ARG:HB2	2.37	0.53
32:DK:38:ILE:HD13	32:DK:61:VAL:HG12	1.91	0.53
32:DK:70:ARG:O	32:DK:71:ARG:HG2	2.08	0.53
32:DK:93:GLN:CG	32:DK:94:PRO:HD2	2.38	0.53
33:DL:79:LEU:CD1	33:DL:112:LEU:HD23	2.39	0.53
35:DN:24:MET:HG2	35:DN:44:LEU:HD13	1.91	0.53
35:DN:42:LYS:CE	35:DN:45:ARG:HG3	2.37	0.53
40:DS:84:ARG:HH21	40:DS:98:LYS:HZ3	1.56	0.53
41:DT:31:VAL:HG22	41:DT:32:LEU:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DU:57:ILE:HD13	42:DU:57:ILE:C	2.29	0.53
45:DY:4:ILE:HG23	45:DY:5:LYS:HD3	1.91	0.53
1:AA:1027:C:H2'	1:AA:1028:C:C6	2.43	0.53
1:AA:600:A:H2'	1:AA:601:G:H8	1.73	0.53
1:AA:633:G:H2'	1:AA:634:C:C6	2.44	0.53
2:AC:40:GLN:O	2:AC:44:LYS:HG3	2.08	0.53
2:AC:85:LYS:O	2:AC:89:VAL:HG23	2.08	0.53
6:AG:129:ASN:HA	6:AG:134:VAL:HG21	1.89	0.53
13:AN:59:GLN:NE2	13:AN:59:GLN:N	2.56	0.53
18:AS:10:ILE:HG22	18:AS:38:THR:H	1.73	0.53
23:BB:1006:C:O2'	23:BB:1007:C:H5'	2.08	0.53
23:BB:1028:A:N6	23:BB:1125:G:H2'	2.24	0.53
23:BB:1684:G:H2'	23:BB:1685:C:C6	2.44	0.53
23:BB:2097:A:H2'	23:BB:2098:U:H6	1.73	0.53
23:BB:910:A:H2	23:BB:2264:C:O2	1.92	0.53
23:BB:2692:G:O2'	23:BB:2693:G:H5'	2.08	0.53
23:BB:2874:C:H2'	23:BB:2875:C:C6	2.43	0.53
23:BB:2899:A:H2'	23:BB:2900:A:H8	1.74	0.53
23:BB:298:G:OP1	42:BU:82:VAL:HA	2.09	0.53
23:BB:460:A:H2'	23:BB:461:C:O4'	2.09	0.53
23:BB:710:U:H2'	23:BB:711:G:C8	2.42	0.53
23:BB:817:C:O2'	23:BB:839:U:H5''	2.09	0.53
23:BB:948:C:H2'	23:BB:949:G:H8	1.74	0.53
25:BC:58:LYS:NZ	25:BC:58:LYS:HB2	2.24	0.53
26:BD:77:ARG:HG2	26:BD:78:GLY:N	2.23	0.53
31:BJ:10:THR:HB	31:BJ:13:ARG:NH1	2.24	0.53
34:BM:108:VAL:O	34:BM:110:GLU:N	2.42	0.53
39:BR:5:PHE:CE2	39:BR:7:SER:HB3	2.44	0.53
40:BS:56:ALA:C	40:BS:58:ALA:H	2.12	0.53
43:BW:76:ARG:CZ	43:BW:76:ARG:HB3	2.38	0.53
1:CA:1026:G:H2'	1:CA:1027:C:C6	2.41	0.53
1:CA:1060:U:H2'	1:CA:1061:G:C8	2.44	0.53
1:CA:1283:U:O2'	1:CA:1284:C:H5'	2.08	0.53
1:CA:140:U:H2'	1:CA:141:G:H8	1.73	0.53
1:CA:149:A:H1'	1:CA:1446:A:C2	2.44	0.53
1:CA:437:U:C2'	1:CA:438:U:O4'	2.51	0.53
1:CA:471:U:H2'	1:CA:472:U:H6	1.73	0.53
1:CA:613:C:H2'	1:CA:614:C:C6	2.44	0.53
1:CA:736:C:H2'	1:CA:737:C:C6	2.43	0.53
20:CB:102:ASN:ND2	20:CB:105:THR:HB	2.24	0.53
4:CE:110:MET:HG2	4:CE:139:THR:OG1	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CE:35:LEU:HD13	4:CE:133:ILE:HA	1.91	0.53
8:CI:51:LEU:HD22	8:CI:56:MET:HG2	1.90	0.53
1:CA:521:G:OP1	11:CL:69:GLU:HB3	2.08	0.53
9:CJ:51:VAL:CG2	13:CN:80:ARG:HB2	2.33	0.53
21:CU:25:ALA:C	21:CU:27:VAL:H	2.10	0.53
48:D1:8:ILE:CD1	48:D1:9:LYS:H	2.22	0.53
22:DA:94:A:O2'	22:DA:95:U:H5'	2.08	0.53
23:DB:1082:U:C4	23:DB:1086:A:N1	2.77	0.53
23:DB:1813:G:N3	25:DC:50:THR:CG2	2.69	0.53
23:DB:1946:U:H2'	23:DB:1947:C:C6	2.43	0.53
23:DB:417:C:H2'	23:DB:418:C:C6	2.44	0.53
23:DB:945:A:H3'	23:DB:946:C:H5''	1.90	0.53
25:DC:152:GLN:HB3	25:DC:153:LEU:HD23	1.90	0.53
23:DB:1791:A:C5'	25:DC:211:ARG:HE	2.20	0.53
25:DC:28:PRO:HD2	25:DC:79:ARG:HE	1.73	0.53
23:DB:2820:A:N6	26:DD:197:THR:HB	2.24	0.53
29:DG:11:PRO:HD2	29:DG:14:VAL:HG21	1.90	0.53
52:DI:1:ALA:HB1	52:DI:2:LYS:HD2	1.90	0.53
52:DI:54:ILE:HD13	52:DI:55:PRO:N	2.22	0.53
23:DB:2641:G:OP1	31:DJ:78:THR:HG22	2.09	0.53
32:DK:108:ARG:HH11	32:DK:108:ARG:HG3	1.73	0.53
23:DB:635:C:P	33:DL:126:ARG:HE	2.32	0.53
33:DL:39:LYS:HZ3	33:DL:39:LYS:N	2.06	0.53
36:DO:30:ARG:HD2	36:DO:97:PHE:HD2	1.74	0.53
37:DP:31:VAL:HG12	37:DP:33:GLU:H	1.72	0.53
39:DR:4:VAL:HG12	39:DR:43:ASN:CB	2.38	0.53
39:DR:44:GLY:N	39:DR:53:PHE:HE2	2.06	0.53
39:DR:78:ARG:O	39:DR:79:ARG:HB2	2.09	0.53
40:DS:27:LYS:HA	40:DS:70:LYS:CB	2.38	0.53
41:DT:7:LEU:HD22	41:DT:7:LEU:H	1.73	0.53
43:DW:21:GLY:HA2	43:DW:25:PHE:CE1	2.43	0.53
44:DX:17:GLU:O	44:DX:22:LEU:HG	2.08	0.53
1:AA:113:G:H2'	1:AA:114:U:C6	2.43	0.53
1:AA:252:U:H2'	1:AA:253:A:H8	1.74	0.53
1:AA:287:U:O2'	1:AA:288:A:H5'	2.08	0.53
2:AC:139:ASN:O	2:AC:143:LEU:HD22	2.09	0.53
9:AJ:35:GLN:H	9:AJ:78:GLU:HB3	1.74	0.53
11:AL:79:ILE:HG22	11:AL:103:CYS:HB2	1.90	0.53
11:AL:80:LEU:HD23	11:AL:97:VAL:HG21	1.90	0.53
13:AN:82:LYS:HE2	13:AN:82:LYS:HA	1.91	0.53
18:AS:31:ARG:HH11	18:AS:31:ARG:HG2	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AS:10:ILE:HD13	18:AS:40:PHE:CE1	2.44	0.53
48:B1:42:VAL:CG1	48:B1:43:ARG:N	2.72	0.53
22:BA:17:C:C2'	22:BA:18:G:H5'	2.39	0.53
23:BB:1060:U:O2	23:BB:1088:A:C8	2.62	0.53
23:BB:1583:A:H4'	23:BB:1585:C:C4	2.44	0.53
23:BB:1671:U:H2'	23:BB:1673:G:OP2	2.08	0.53
23:BB:1864:U:O2'	23:BB:1865:U:H5'	2.08	0.53
23:BB:257:C:H2'	23:BB:258:G:O4'	2.09	0.53
23:BB:2634:A:H2'	23:BB:2635:A:C8	2.44	0.53
23:BB:309:A:N3	23:BB:329:G:O2'	2.42	0.53
23:BB:560:C:H3'	23:BB:561:G:C8	2.44	0.53
23:BB:664:G:H2'	23:BB:665:U:C6	2.44	0.53
23:BB:847:U:O4'	23:BB:847:U:O2	2.27	0.53
26:BD:131:ASP:OD2	26:BD:133:THR:HG22	2.08	0.53
26:BD:60:VAL:HB	26:BD:63:PRO:CG	2.38	0.53
27:BE:44:ARG:N	27:BE:44:ARG:HD3	2.24	0.53
29:BG:157:LYS:CG	29:BG:159:LYS:HZ2	2.21	0.53
31:BJ:78:THR:O	31:BJ:80:HIS:N	2.42	0.53
31:BJ:72:LYS:CB	31:BJ:90:GLU:HB2	2.34	0.53
45:BY:4:ILE:HG12	45:BY:5:LYS:HE2	1.90	0.53
46:BZ:5:ILE:HD13	46:BZ:51:VAL:HG12	1.91	0.53
1:CA:276:G:H5'	16:CQ:16:MET:SD	2.48	0.53
1:CA:817:C:H1'	1:CA:819:A:C5'	2.38	0.53
1:CA:952:U:H2'	1:CA:953:G:C8	2.41	0.53
20:CB:40:ILE:HG21	20:CB:200:PRO:O	2.09	0.53
2:CC:49:ALA:HB2	2:CC:74:ILE:HG21	1.91	0.53
2:CC:89:VAL:HG23	2:CC:90:VAL:N	2.23	0.53
3:CD:104:MET:SD	3:CD:179:GLY:HA2	2.47	0.53
4:CE:136:VAL:HG23	4:CE:137:ARG:N	2.23	0.53
12:CM:15:VAL:HG13	12:CM:40:GLU:HB2	1.89	0.53
15:CP:3:THR:HB	15:CP:66:THR:O	2.07	0.53
49:D2:6:GLN:HB3	49:D2:7:PRO:HD2	1.89	0.53
51:D4:16:ILE:C	51:D4:18:LYS:H	2.12	0.53
23:DB:1341:G:H2'	23:DB:1397:U:O2'	2.09	0.53
23:DB:1583:A:H4'	23:DB:1585:C:C4	2.44	0.53
23:DB:1733:G:H2'	23:DB:1734:G:H8	1.73	0.53
23:DB:1826:G:H2'	23:DB:1827:U:C6	2.44	0.53
23:DB:2348:U:O2'	23:DB:2349:G:H5'	2.09	0.53
23:DB:2718:G:O2'	23:DB:2719:G:H5'	2.09	0.53
23:DB:633:A:C2'	23:DB:634:C:H5'	2.37	0.53
23:DB:937:C:H2'	23:DB:938:G:C8	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:123:ILE:HG12	25:DC:135:PRO:CD	2.39	0.53
25:DC:224:MET:O	25:DC:225:ASN:HB2	2.07	0.53
27:DE:15:SER:O	27:DE:17:THR:HG22	2.09	0.53
27:DE:161:ALA:O	27:DE:162:ARG:HB2	2.07	0.53
31:DJ:40:HIS:O	38:DQ:66:ALA:HB1	2.09	0.53
33:DL:58:TYR:HB3	50:D3:13:PHE:HZ	1.73	0.53
35:DN:32:GLU:O	35:DN:114:GLU:HA	2.08	0.53
36:DO:25:ARG:O	36:DO:26:LEU:HD12	2.09	0.53
38:DQ:40:LYS:HB2	38:DQ:40:LYS:NZ	2.23	0.53
40:DS:31:GLN:O	40:DS:35:ILE:HG13	2.08	0.53
1:AA:613:C:H2'	1:AA:614:C:C6	2.44	0.53
1:AA:640:A:O2'	1:AA:641:U:H5'	2.08	0.53
1:AA:672:U:H2'	1:AA:673:A:C8	2.43	0.53
1:AA:711:G:O2'	1:AA:712:A:H5'	2.09	0.53
7:AH:63:LYS:HG2	7:AH:70:VAL:HG21	1.90	0.53
9:AJ:87:LEU:HB3	9:AJ:88:MET:HE3	1.89	0.53
49:B2:25:LYS:H	49:B2:25:LYS:CE	2.21	0.53
23:BB:136:G:H2'	23:BB:137:U:C5	2.43	0.53
23:BB:1745:A:O2'	23:BB:1746:A:H5'	2.09	0.53
23:BB:1796:U:H2'	23:BB:1797:G:H8	1.73	0.53
23:BB:2229:U:H2'	23:BB:2230:G:H8	1.74	0.53
23:BB:197:A:N6	23:BB:2430:A:H2'	2.24	0.53
23:BB:2628:C:O2'	23:BB:2781:A:H2'	2.09	0.53
23:BB:340:A:O2'	23:BB:341:C:H5'	2.08	0.53
23:BB:633:A:O5'	23:BB:633:A:C8	2.57	0.53
23:BB:737:C:O2'	23:BB:738:G:H5'	2.09	0.53
25:BC:16:VAL:O	25:BC:16:VAL:HG12	2.08	0.53
27:BE:108:ILE:HD12	27:BE:108:ILE:H	1.73	0.53
29:BG:100:ASN:O	29:BG:115:GLN:HA	2.09	0.53
29:BG:29:ASN:N	29:BG:29:ASN:HD22	2.05	0.53
37:BP:70:GLU:O	37:BP:71:ARG:HB2	2.08	0.53
39:BR:39:LEU:HD12	39:BR:60:LYS:HB2	1.90	0.53
42:BU:39:ASN:HD21	42:BU:42:LYS:HE2	1.71	0.53
42:BU:59:GLU:O	42:BU:60:LYS:HB2	2.08	0.53
42:BU:60:LYS:CG	42:BU:61:GLU:H	2.22	0.53
42:BU:71:ILE:HD13	42:BU:71:ILE:C	2.29	0.53
24:BV:30:ILE:HD11	24:BV:63:ILE:HD12	1.90	0.53
1:CA:1158:C:O3'	20:CB:131:LYS:HD3	2.08	0.53
1:CA:1277:C:H1'	1:CA:1282:C:O2	2.09	0.53
1:CA:1313:U:OP2	18:CS:5:LYS:HA	2.09	0.53
1:CA:764:C:C2'	1:CA:765:G:H5'	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:122:GLN:HB3	2:CC:127:VAL:HG11	1.89	0.53
2:CC:166:TRP:HE1	2:CC:168:ARG:HB2	1.73	0.53
4:CE:33:THR:HB	4:CE:49:TYR:CE1	2.43	0.53
5:CF:9:MET:HB2	5:CF:57:ALA:HB1	1.90	0.53
7:CH:71:VAL:HG23	7:CH:71:VAL:O	2.09	0.53
16:CQ:16:MET:HB2	16:CQ:19:SER:HB2	1.91	0.53
1:CA:1311:A:N6	18:CS:1:PRO:HD3	2.21	0.53
22:DA:10:G:H2'	22:DA:11:C:O4'	2.09	0.53
23:DB:1138:G:H2'	23:DB:1139:G:O4'	2.09	0.53
23:DB:1571:A:H2'	23:DB:1572:A:H8	1.73	0.53
23:DB:1684:G:H2'	23:DB:1685:C:C6	2.44	0.53
23:DB:1857:G:HO2'	23:DB:1858:A:H8	1.53	0.53
23:DB:2241:A:H2'	23:DB:2242:G:H8	1.72	0.53
23:DB:2467:C:OP2	51:D4:7:VAL:HG12	2.09	0.53
23:DB:2529:G:H5'	29:DG:175:LYS:HB3	1.89	0.53
23:DB:2628:C:O2'	23:DB:2781:A:H2'	2.09	0.53
23:DB:521:U:H2'	23:DB:522:A:C8	2.43	0.53
25:DC:268:ARG:O	25:DC:269:ARG:HB2	2.09	0.53
25:DC:53:ILE:CD1	25:DC:218:THR:HA	2.39	0.53
26:DD:154:LYS:C	26:DD:156:PHE:H	2.12	0.53
26:DD:21:SER:C	26:DD:23:PRO:HD3	2.29	0.53
27:DE:189:THR:HG23	27:DE:194:LYS:HB2	1.89	0.53
29:DG:3:VAL:O	29:DG:4:ALA:HB2	2.08	0.53
30:DH:94:ILE:HG23	30:DH:98:ASP:HB2	1.90	0.53
32:DK:41:ILE:HG13	32:DK:42:THR:H	1.74	0.53
33:DL:109:LYS:HG2	33:DL:126:ARG:CD	2.38	0.53
34:DM:33:LEU:HG	34:DM:113:ALA:HB1	1.91	0.53
35:DN:24:MET:CE	35:DN:36:THR:HG21	2.39	0.53
35:DN:29:VAL:HG21	35:DN:75:ILE:HB	1.90	0.53
35:DN:107:ASN:ND2	40:DS:40:ASN:HB3	2.23	0.53
24:DV:63:ILE:N	24:DV:63:ILE:HD12	2.24	0.53
44:DX:7:ARG:NH1	44:DX:7:ARG:HB3	2.24	0.53
1:AA:44:A:O2'	1:AA:45:G:H5'	2.09	0.53
1:AA:476:U:H2'	1:AA:477:C:H6	1.73	0.53
1:AA:963:G:H21	9:AJ:56:HIS:HE1	1.55	0.53
10:AK:105:ARG:NH2	21:AU:10:PRO:HG3	2.24	0.53
11:AL:28:GLN:HG3	11:AL:80:LEU:HD21	1.90	0.53
13:AN:51:PRO:HB2	13:AN:54:SER:HB3	1.91	0.53
14:AO:57:ARG:HD3	14:AO:57:ARG:C	2.30	0.53
51:B4:14:CYS:SG	51:B4:27:CYS:HB3	2.48	0.53
23:BB:1549:A:H2'	23:BB:1550:C:C6	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1826:G:H2'	23:BB:1827:U:C6	2.44	0.53
23:BB:1846:G:H2'	23:BB:1847:A:O4'	2.09	0.53
23:BB:184:C:H2'	23:BB:185:G:H8	1.72	0.53
23:BB:2213:U:O2	23:BB:2213:U:H2'	2.07	0.53
23:BB:2263:C:H2'	23:BB:2264:C:H6	1.74	0.53
23:BB:2308:G:N3	23:BB:2308:G:H5''	2.24	0.53
23:BB:337:C:H2'	23:BB:338:G:O4'	2.09	0.53
26:BD:109:VAL:HG11	26:BD:193:VAL:HG23	1.90	0.53
29:BG:163:TYR:HB2	29:BG:166:GLU:HB2	1.90	0.53
31:BJ:101:ILE:HG23	31:BJ:102:GLU:OE1	2.08	0.53
31:BJ:56:VAL:C	31:BJ:57:LEU:HD22	2.29	0.53
34:BM:114:ARG:HA	34:BM:117:PHE:HB3	1.91	0.53
23:BB:2820:A:O5'	35:BN:4:ARG:HA	2.08	0.53
35:BN:57:THR:HG23	35:BN:58:ASP:H	1.74	0.53
38:BQ:92:LYS:HE3	38:BQ:93:ILE:HB	1.91	0.53
44:BX:14:LEU:O	44:BX:17:GLU:HB2	2.09	0.53
1:CA:462:G:H2'	1:CA:463:U:C6	2.44	0.53
1:CA:975:A:C2'	1:CA:976:G:OP2	2.57	0.53
20:CB:209:VAL:C	20:CB:213:LEU:HD12	2.29	0.53
2:CC:63:ILE:HG12	2:CC:65:VAL:HG23	1.91	0.53
3:CD:170:LEU:HA	3:CD:182:LYS:HG3	1.91	0.53
11:CL:97:VAL:O	11:CL:97:VAL:HG23	2.09	0.53
23:DB:2886:A:C5	47:D0:27:LEU:HG	2.43	0.53
23:DB:1105:U:H2'	23:DB:1106:G:C8	2.41	0.53
23:DB:1430:G:H2'	23:DB:1431:A:H8	1.74	0.53
23:DB:18:U:H2'	23:DB:19:A:H8	1.74	0.53
23:DB:2028:U:H2'	23:DB:2029:G:C8	2.44	0.53
23:DB:443:A:H5''	23:DB:444:C:OP1	2.09	0.53
23:DB:544:C:H2'	23:DB:545:U:N3	2.24	0.53
23:DB:766:U:H2'	23:DB:767:U:C6	2.44	0.53
23:DB:934:U:H2'	23:DB:935:C:C6	2.44	0.53
25:DC:124:LYS:HB2	25:DC:125:PRO:CD	2.39	0.53
25:DC:181:ARG:O	25:DC:181:ARG:HG3	2.09	0.53
25:DC:231:HIS:HD1	25:DC:242:HIS:HA	1.74	0.53
25:DC:32:LEU:N	25:DC:32:LEU:HD23	2.23	0.53
28:DF:177:ARG:CZ	28:DF:178:LYS:HA	2.39	0.53
28:DF:39:VAL:HG13	28:DF:84:ILE:CG1	2.39	0.53
23:DB:632:A:O3'	33:DL:69:ARG:HD3	2.09	0.53
24:DV:1:MET:SD	24:DV:62:THR:HG23	2.49	0.53
43:DW:42:THR:N	43:DW:65:LYS:HG2	2.23	0.53
45:DY:50:VAL:CG1	45:DY:53:MET:HG2	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DZ:41:HIS:O	46:DZ:43:PHE:HD1	1.92	0.53
1:AA:1033:G:H3'	1:AA:1034:G:H5''	1.91	0.53
1:AA:1392:G:H2'	1:AA:1393:U:C6	2.44	0.53
1:AA:588:G:H5'	7:AH:2:MET:O	2.09	0.53
1:AA:777:A:H2'	1:AA:778:G:C8	2.43	0.53
2:AC:119:ILE:O	2:AC:123:LEU:HG	2.09	0.53
2:AC:90:VAL:HG23	2:AC:98:ALA:HB3	1.90	0.53
5:AF:3:HIS:HB2	5:AF:92:THR:OG1	2.09	0.53
6:AG:143:MET:O	6:AG:147:ASN:HB2	2.09	0.53
16:AQ:45:VAL:HG11	16:AQ:74:LEU:HB2	1.90	0.53
16:AQ:68:LYS:O	16:AQ:69:THR:HB	2.09	0.53
16:AQ:57:VAL:HB	16:AQ:79:GLU:H	1.74	0.53
22:BA:78:A:H2'	22:BA:79:G:O4'	2.09	0.53
23:BB:2688:G:H1'	23:BB:2721:A:H61	1.74	0.53
23:BB:2898:U:H2'	23:BB:2899:A:C8	2.44	0.53
23:BB:503:A:C3'	23:BB:504:A:H5'	2.39	0.53
28:BF:4:HIS:O	28:BF:7:TYR:HB3	2.09	0.53
29:BG:102:ILE:O	29:BG:113:ASP:HA	2.08	0.53
29:BG:25:ILE:O	29:BG:31:GLU:HA	2.09	0.53
33:BL:14:LYS:CA	33:BL:14:LYS:HZ3	2.08	0.53
34:BM:65:ILE:HA	34:BM:101:VAL:HA	1.90	0.53
23:BB:2484:G:O2'	34:BM:45:GLN:NE2	2.41	0.53
35:BN:118:ARG:O	35:BN:118:ARG:HG3	2.09	0.53
36:BO:62:LEU:HD12	36:BO:64:TYR:HB2	1.90	0.53
40:BS:20:VAL:O	40:BS:23:LEU:HB2	2.09	0.53
40:BS:86:MET:SD	40:BS:96:ILE:HD11	2.49	0.53
43:BW:16:GLU:O	43:BW:18:LYS:HD2	2.09	0.53
1:CA:1478:U:H2'	1:CA:1479:C:C6	2.43	0.53
1:CA:203:G:N2	1:CA:205:A:N6	2.56	0.53
1:CA:556:C:O2'	1:CA:557:G:H5'	2.08	0.53
1:CA:640:A:O2'	1:CA:641:U:H5'	2.08	0.53
1:CA:858:G:O6	1:CA:869:G:H3'	2.09	0.53
1:CA:883:C:O2'	1:CA:884:U:H5'	2.09	0.53
20:CB:150:ILE:O	20:CB:150:ILE:HG12	2.09	0.53
20:CB:16:GLY:CA	20:CB:40:ILE:H	2.21	0.53
20:CB:53:LEU:HD13	20:CB:212:TYR:OH	2.09	0.53
2:CC:53:ARG:HA	2:CC:113:LYS:NZ	2.23	0.53
2:CC:23:ALA:HB1	2:CC:27:GLU:HB2	1.91	0.53
2:CC:91:ALA:HB2	2:CC:98:ALA:HB3	1.90	0.53
15:CP:44:SER:CB	15:CP:46:LYS:HG2	2.35	0.53
47:D0:51:ARG:O	47:D0:51:ARG:HD3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:250:G:C5'	50:D3:7:ARG:HG2	2.35	0.53
23:DB:1368:G:OP1	49:D2:25:LYS:HG3	2.09	0.53
23:DB:1594:U:H2'	23:DB:1595:C:H6	1.74	0.53
23:DB:1846:G:H2'	23:DB:1847:A:O4'	2.09	0.53
23:DB:228:C:H4'	23:DB:229:C:H5''	1.90	0.53
23:DB:2489:U:O2'	23:DB:2490:G:H5'	2.08	0.53
23:DB:2512:C:H2'	23:DB:2513:A:O4'	2.08	0.53
23:DB:2634:A:H2'	23:DB:2635:A:C8	2.44	0.53
26:DD:149:ASN:H	26:DD:151:THR:HG23	1.73	0.53
23:DB:2313:C:H5'	28:DF:36:ASN:ND2	2.23	0.53
31:DJ:109:LEU:HD22	31:DJ:115:GLY:O	2.09	0.53
37:DP:89:GLY:HA2	37:DP:112:ARG:NH2	2.23	0.53
24:DV:53:LYS:NZ	24:DV:55:GLU:HG2	2.23	0.53
1:AA:537:G:H2'	1:AA:538:G:H8	1.74	0.52
20:AB:30:ILE:HG22	20:AB:32:GLY:H	1.74	0.52
4:AE:156:ARG:HB2	7:AH:43:GLY:HA3	1.91	0.52
6:AG:130:LYS:N	6:AG:134:VAL:HG11	2.23	0.52
8:AI:115:VAL:HG21	9:AJ:62:ARG:HB2	1.90	0.52
12:AM:27:THR:O	12:AM:30:LYS:HB3	2.09	0.52
19:AT:28:ARG:HA	19:AT:31:ILE:HD12	1.91	0.52
48:B1:42:VAL:O	48:B1:43:ARG:HB2	2.09	0.52
23:BB:1050:A:H2'	23:BB:1051:G:O4'	2.10	0.52
23:BB:1183:U:H2'	23:BB:1184:U:C6	2.43	0.52
23:BB:1844:C:OP1	25:BC:252:LYS:HA	2.09	0.52
23:BB:2144:G:N2	23:BB:2148:G:N3	2.57	0.52
23:BB:2218:G:H2'	23:BB:2219:U:H6	1.73	0.52
23:BB:2305:U:H2'	23:BB:2306:C:C6	2.43	0.52
23:BB:2675:A:N1	23:BB:2732:G:O6	2.42	0.52
23:BB:877:A:N1	23:BB:900:A:N7	2.57	0.52
27:BE:144:GLU:HG3	27:BE:145:ASP:H	1.73	0.52
28:BF:106:ALA:HB1	28:BF:136:ILE:HG23	1.90	0.52
33:BL:106:GLU:HG2	33:BL:107:PHE:H	1.74	0.52
33:BL:126:ARG:HH21	33:BL:131:ALA:HB2	1.73	0.52
33:BL:56:PRO:O	33:BL:59:ARG:HB2	2.08	0.52
34:BM:15:GLY:O	34:BM:16:ARG:HB3	2.09	0.52
34:BM:7:THR:O	34:BM:8:LYS:HD3	2.09	0.52
38:BQ:84:LYS:N	38:BQ:84:LYS:HE3	2.24	0.52
39:BR:63:VAL:HB	39:BR:101:ILE:HG12	1.91	0.52
38:BQ:88:GLU:HG2	39:BR:53:PHE:CD2	2.45	0.52
40:BS:11:ARG:HD3	40:BS:11:ARG:O	2.08	0.52
44:BX:29:ARG:CA	44:BX:34:SER:HB2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1142:G:H2'	1:CA:1143:G:C8	2.44	0.52
1:CA:113:G:H2'	1:CA:114:U:C6	2.44	0.52
1:CA:1171:A:H2'	1:CA:1172:C:C6	2.44	0.52
1:CA:1244:G:H2'	1:CA:1245:C:H6	1.71	0.52
2:CC:151:GLU:HA	2:CC:166:TRP:HA	1.90	0.52
5:CF:25:TYR:O	5:CF:29:ILE:HG13	2.09	0.52
4:CE:156:ARG:HD2	7:CH:42:GLU:O	2.09	0.52
13:CN:43:ALA:O	13:CN:47:LEU:HB2	2.09	0.52
16:CQ:60:ILE:HD13	16:CQ:60:ILE:H	1.74	0.52
51:D4:19:ARG:HB3	51:D4:19:ARG:HH11	1.73	0.52
23:DB:1042:G:H2'	23:DB:1043:C:C6	2.45	0.52
23:DB:1060:U:O2	23:DB:1088:A:C8	2.63	0.52
23:DB:1186:G:H2'	23:DB:1187:G:O4'	2.09	0.52
23:DB:1864:U:O2'	23:DB:1865:U:H5'	2.08	0.52
23:DB:2047:C:H2'	23:DB:2048:G:H8	1.74	0.52
23:DB:204:A:H4'	23:DB:205:G:OP1	2.09	0.52
25:DC:37:SER:H	25:DC:62:ARG:CB	2.21	0.52
27:DE:120:VAL:HG12	27:DE:121:VAL:N	2.18	0.52
28:DF:39:VAL:HG13	28:DF:84:ILE:HB	1.91	0.52
33:DL:69:ARG:HG3	33:DL:71:ALA:O	2.09	0.52
33:DL:90:VAL:CG1	33:DL:122:VAL:HG22	2.39	0.52
33:DL:92:LEU:HA	33:DL:96:LYS:HG2	1.91	0.52
23:DB:929:U:H1'	45:DY:24:LEU:O	2.09	0.52
1:AA:1115:U:H2'	1:AA:1116:U:C6	2.45	0.52
1:AA:1438:G:O2'	1:AA:1439:G:H5'	2.08	0.52
20:AB:160:LEU:O	20:AB:182:VAL:HA	2.09	0.52
6:AG:16:LYS:HG3	6:AG:43:TYR:OH	2.08	0.52
9:AJ:5:ARG:HD3	9:AJ:5:ARG:N	2.24	0.52
13:AN:81:ILE:O	13:AN:84:ARG:HB3	2.09	0.52
16:AQ:45:VAL:HG12	16:AQ:46:HIS:N	2.21	0.52
23:BB:1151:A:H2'	23:BB:1152:C:H6	1.74	0.52
23:BB:154:U:H2'	23:BB:155:A:H8	1.74	0.52
23:BB:2213:U:H2'	23:BB:2214:C:H5'	1.92	0.52
23:BB:2348:U:O2'	23:BB:2349:G:H5'	2.09	0.52
23:BB:2649:C:H2'	23:BB:2650:U:H6	1.73	0.52
23:BB:1664:A:H1'	23:BB:2726:A:N1	2.23	0.52
23:BB:279:A:N1	23:BB:362:A:H5'	2.24	0.52
23:BB:328:U:H4'	42:BU:65:GLN:HE21	1.74	0.52
23:BB:49:A:H4'	23:BB:49:A:OP1	2.09	0.52
25:BC:81:GLU:HB3	25:BC:90:ILE:CG2	2.40	0.52
26:BD:15:PHE:HB3	26:BD:18:ASP:HB3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2750:A:H3'	29:BG:3:VAL:HG13	1.90	0.52
30:BH:41:LYS:H	30:BH:41:LYS:HD2	1.74	0.52
52:BI:81:LYS:HG3	52:BI:82:ALA:N	2.24	0.52
31:BJ:114:LEU:O	31:BJ:117:ALA:HB3	2.10	0.52
34:BM:32:GLY:HA2	34:BM:101:VAL:O	2.09	0.52
37:BP:47:ILE:HG23	37:BP:63:ILE:HG23	1.91	0.52
23:BB:2849:U:O4	37:BP:96:LEU:HD21	2.09	0.52
38:BQ:50:ARG:HA	38:BQ:53:LYS:CE	2.39	0.52
41:BT:50:LEU:C	41:BT:52:GLU:H	2.13	0.52
41:BT:3:ARG:HG2	41:BT:7:LEU:HB3	1.92	0.52
1:CA:1337:G:H5''	1:CA:1338:G:OP1	2.09	0.52
1:CA:1480:A:H2'	1:CA:1481:U:C6	2.43	0.52
1:CA:685:G:O2'	1:CA:686:U:H5'	2.09	0.52
20:CB:23:ASN:HD22	20:CB:24:PRO:CD	2.22	0.52
5:CF:85:ILE:HG22	5:CF:86:ARG:H	1.73	0.52
5:CF:38:ARG:NH1	5:CF:98:GLU:H	1.98	0.52
9:CJ:41:PRO:HA	9:CJ:72:ARG:CD	2.38	0.52
9:CJ:5:ARG:O	9:CJ:5:ARG:HG2	2.09	0.52
13:CN:5:MET:HA	13:CN:8:ARG:HG3	1.92	0.52
16:CQ:10:ARG:CZ	16:CQ:11:VAL:H	2.22	0.52
18:CS:43:MET:HB3	18:CS:61:VAL:HG11	1.91	0.52
48:D1:10:LEU:HD13	48:D1:50:GLU:OE1	2.10	0.52
50:D3:33:THR:CG2	50:D3:40:LYS:HD2	2.38	0.52
23:DB:2359:C:H4'	50:D3:51:LYS:HD2	1.91	0.52
22:DA:87:U:O2'	22:DA:88:C:H2'	2.08	0.52
23:DB:178:G:O2'	23:DB:179:C:H5'	2.09	0.52
23:DB:1911:U:H2'	23:DB:1918:A:N1	2.22	0.52
23:DB:323:C:H5	27:DE:165:HIS:HB2	1.74	0.52
30:DH:135:HIS:CD2	30:DH:137:GLU:H	2.26	0.52
37:DP:91:VAL:HG23	37:DP:113:LEU:HB3	1.92	0.52
37:DP:7:LEU:CD2	37:DP:11:GLN:HE21	2.22	0.52
37:DP:77:SER:HB2	37:DP:78:PRO:HD3	1.90	0.52
23:DB:1754:A:OP1	37:DP:95:LYS:HB2	2.10	0.52
38:DQ:26:ALA:CA	38:DQ:30:VAL:HG23	2.39	0.52
23:DB:997:G:OP1	38:DQ:92:LYS:HB3	2.09	0.52
1:AA:214:C:H2'	1:AA:215:C:H6	1.72	0.52
1:AA:333:U:H2'	1:AA:334:C:H6	1.74	0.52
1:AA:392:C:H2'	1:AA:393:A:H8	1.75	0.52
1:AA:556:C:O2'	1:AA:557:G:H5'	2.08	0.52
1:AA:858:G:O6	1:AA:869:G:H3'	2.09	0.52
2:AC:71:ARG:HB3	2:AC:74:ILE:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:76:ILE:C	2:AC:82:ASP:HB2	2.30	0.52
2:AC:78:LYS:HG3	2:AC:78:LYS:O	2.10	0.52
3:AD:117:VAL:HG12	3:AD:130:ASN:C	2.30	0.52
6:AG:16:LYS:HD3	6:AG:16:LYS:C	2.29	0.52
8:AI:48:ARG:HA	8:AI:51:LEU:HD13	1.92	0.52
8:AI:47:VAL:HG12	8:AI:78:ILE:HB	1.91	0.52
1:AA:552:U:H4'	11:AL:82:ARG:HG2	1.92	0.52
9:AJ:56:HIS:N	13:AN:80:ARG:HH22	2.07	0.52
23:BB:1010:A:H5'	38:BQ:61:ILE:HD12	1.91	0.52
23:BB:2028:U:H2'	23:BB:2029:G:C8	2.44	0.52
23:BB:2618:G:O2'	26:BD:154:LYS:O	2.26	0.52
23:BB:2719:G:H4'	23:BB:2846:G:O3'	2.09	0.52
23:BB:522:A:H2'	23:BB:523:C:C6	2.45	0.52
23:BB:600:G:H2'	23:BB:601:C:C6	2.45	0.52
23:BB:995:C:N4	31:BJ:2:LYS:HB3	2.25	0.52
25:BC:114:GLN:HB2	25:BC:124:LYS:NZ	2.24	0.52
25:BC:46:GLY:O	25:BC:47:ARG:CB	2.56	0.52
25:BC:53:ILE:HD13	25:BC:218:THR:CG2	2.30	0.52
23:BB:2579:C:O2'	26:BD:138:LEU:HD13	2.10	0.52
31:BJ:34:ARG:HB3	31:BJ:39:LYS:HD2	1.91	0.52
32:BK:41:ILE:HG23	32:BK:42:THR:N	2.24	0.52
39:BR:90:ARG:NH1	39:BR:91:GLN:H	2.08	0.52
24:BV:14:LYS:HZ2	24:BV:18:ARG:HD2	1.72	0.52
1:CA:1000:A:H2'	1:CA:1001:C:H6	1.71	0.52
1:CA:1049:U:O2'	1:CA:1050:G:OP2	2.26	0.52
1:CA:1057:G:H5''	2:CC:153:SER:HB2	1.91	0.52
1:CA:1330:U:OP1	12:CM:22:TYR:O	2.27	0.52
1:CA:394:G:H2'	1:CA:395:C:H6	1.75	0.52
1:CA:537:G:H2'	1:CA:538:G:H8	1.75	0.52
1:CA:663:A:O2'	1:CA:664:G:H5'	2.10	0.52
3:CD:171:GLU:HG3	3:CD:182:LYS:CD	2.39	0.52
4:CE:35:LEU:CD1	4:CE:133:ILE:HA	2.40	0.52
5:CF:38:ARG:NH2	5:CF:63:ASN:HD21	2.07	0.52
5:CF:3:HIS:NE2	5:CF:95:ALA:HB2	2.24	0.52
6:CG:145:GLU:H	6:CG:148:LYS:HB2	1.75	0.52
6:CG:77:ARG:HD3	6:CG:79:VAL:HG23	1.92	0.52
1:CA:1060:U:C4'	9:CJ:54:SER:HB2	2.39	0.52
10:CK:109:ILE:H	10:CK:109:ILE:HD12	1.74	0.52
13:CN:31:SER:O	13:CN:40:ARG:HD3	2.09	0.52
1:CA:625:U:H4'	15:CP:16:PHE:CE2	2.44	0.52
23:DB:1034:G:H2'	23:DB:1035:U:C6	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:136:G:C2	41:DT:3:ARG:NH2	2.77	0.52
23:DB:1374:G:H2'	23:DB:1375:U:C6	2.45	0.52
23:DB:1733:G:H2'	23:DB:1734:G:C8	2.45	0.52
23:DB:18:U:H2'	23:DB:19:A:C8	2.44	0.52
23:DB:19:A:H2'	23:DB:20:C:C6	2.44	0.52
23:DB:19:A:H2'	23:DB:20:C:H6	1.75	0.52
23:DB:20:C:O2'	23:DB:21:A:H5'	2.09	0.52
23:DB:2675:A:N1	23:DB:2732:G:O6	2.42	0.52
23:DB:2734:A:H2'	23:DB:2735:G:H5'	1.91	0.52
23:DB:324:A:N6	23:DB:339:U:H5'	2.24	0.52
23:DB:664:G:H2'	23:DB:665:U:C6	2.43	0.52
23:DB:721:A:H2'	23:DB:722:A:H8	1.74	0.52
32:DK:71:ARG:CB	32:DK:72:PRO:CD	2.87	0.52
33:DL:108:ALA:O	33:DL:109:LYS:HB2	2.09	0.52
37:DP:52:ARG:CB	37:DP:60:VAL:HG11	2.36	0.52
37:DP:6:GLN:N	37:DP:6:GLN:CD	2.63	0.52
39:DR:26:ASP:O	39:DR:27:ILE:HB	2.09	0.52
42:DU:21:ARG:HG3	42:DU:21:ARG:NH1	2.23	0.52
1:AA:1225:A:H3'	1:AA:1226:C:H6	1.74	0.52
1:AA:1390:U:H2'	1:AA:1391:U:C6	2.44	0.52
1:AA:317:U:H2'	1:AA:318:G:H8	1.73	0.52
20:AB:163:ILE:HD11	20:AB:209:VAL:HG12	1.92	0.52
5:AF:68:GLN:O	5:AF:71:ILE:HG13	2.09	0.52
6:AG:13:PRO:HA	6:AG:23:ALA:HB2	1.90	0.52
7:AH:64:TYR:CB	7:AH:69:ALA:HA	2.38	0.52
8:AI:61:ASP:O	8:AI:62:LEU:HD13	2.09	0.52
17:AR:31:TYR:CD2	17:AR:54:LEU:HD21	2.45	0.52
22:BA:54:G:H2'	22:BA:55:U:C6	2.44	0.52
23:BB:1242:U:H2'	23:BB:1243:C:C6	2.44	0.52
23:BB:1739:A:H2'	23:BB:1740:G:O4'	2.10	0.52
23:BB:175:G:H2'	23:BB:176:A:C8	2.45	0.52
23:BB:2553:G:H2'	23:BB:2554:U:H4'	1.90	0.52
23:BB:2734:A:H2'	23:BB:2735:G:H5'	1.91	0.52
23:BB:2741:A:H2'	23:BB:2742:G:O4'	2.08	0.52
23:BB:766:U:H2'	23:BB:767:U:C6	2.45	0.52
25:BC:49:THR:O	25:BC:50:THR:HB	2.10	0.52
32:BK:101:GLY:O	32:BK:119:ALA:HB1	2.08	0.52
33:BL:30:THR:HA	33:BL:36:LYS:CE	2.37	0.52
38:BQ:99:VAL:HG23	38:BQ:100:PHE:HD1	1.73	0.52
23:BB:143:C:O2'	41:BT:5:GLU:HA	2.09	0.52
42:BU:100:GLU:O	42:BU:102:ILE:HG13	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:52:ASN:C	42:BU:54:PRO:HD3	2.30	0.52
23:BB:2386:A:H4'	43:BW:38:ARG:HB2	1.91	0.52
43:BW:36:ILE:HB	43:BW:68:PHE:CD2	2.41	0.52
1:CA:1302:C:H5''	1:CA:1303:C:OP2	2.10	0.52
4:CE:80:LEU:HD23	4:CE:122:VAL:HG11	1.91	0.52
13:CN:60:ARG:O	13:CN:61:ASN:HB2	2.09	0.52
14:CO:61:GLN:O	14:CO:65:LEU:HG	2.08	0.52
48:D1:20:TYR:O	48:D1:21:THR:HB	2.09	0.52
48:D1:46:VAL:HG22	48:D1:47:ILE:N	2.24	0.52
49:D2:12:ARG:HH21	49:D2:16:HIS:CB	2.22	0.52
23:DB:1099:G:N7	52:DI:3:LYS:HD3	2.25	0.52
23:DB:2280:G:O6	43:DW:10:ARG:NH2	2.41	0.52
23:DB:265:A:H62	23:DB:427:U:HO2'	1.56	0.52
23:DB:2774:C:H2'	23:DB:2775:G:O4'	2.09	0.52
23:DB:2794:C:H2'	23:DB:2795:C:C6	2.44	0.52
23:DB:288:U:O2'	23:DB:289:G:H5'	2.08	0.52
23:DB:2898:U:H2'	23:DB:2899:A:C8	2.45	0.52
23:DB:448:U:H3'	27:DE:79:ARG:NE	2.23	0.52
23:DB:533:G:H5'	38:DQ:23:TYR:HD1	1.73	0.52
25:DC:226:PRO:HG3	25:DC:232:GLY:O	2.10	0.52
25:DC:22:GLU:CD	25:DC:202:ARG:HE	2.12	0.52
35:DN:10:LEU:HG	35:DN:11:ASN:N	2.24	0.52
35:DN:4:ARG:N	35:DN:4:ARG:CZ	2.73	0.52
38:DQ:52:ARG:HG2	38:DQ:53:LYS:NZ	2.24	0.52
24:DV:9:ARG:HE	24:DV:20:LEU:HD11	1.74	0.52
1:AA:1170:A:O5'	1:AA:1170:A:H8	1.92	0.52
1:AA:1256:A:O2'	1:AA:1257:A:H5''	2.09	0.52
1:AA:246:A:C2	1:AA:282:A:C5	2.98	0.52
1:AA:303:A:H2'	1:AA:304:U:O4'	2.10	0.52
1:AA:33:A:H2'	1:AA:34:C:C6	2.45	0.52
1:AA:520:A:OP2	11:AL:47:ALA:HB1	2.09	0.52
1:AA:860:A:H2'	1:AA:861:G:O4'	2.09	0.52
20:AB:95:TRP:HH2	20:AB:100:LEU:HB2	1.74	0.52
20:AB:71:THR:HB	20:AB:167:HIS:CD2	2.44	0.52
20:AB:172:ILE:HG22	20:AB:176:ASN:ND2	2.25	0.52
20:AB:20:ARG:HG3	20:AB:20:ARG:HH11	1.73	0.52
20:AB:36:LYS:H	20:AB:36:LYS:HE3	1.71	0.52
8:AI:29:ILE:HD11	8:AI:37:TYR:HB3	1.92	0.52
17:AR:20:ILE:HG23	17:AR:53:GLN:NE2	2.25	0.52
23:BB:1000:A:H2'	23:BB:1001:A:C8	2.45	0.52
23:BB:142:A:C4	41:BT:2:ILE:HD12	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1506:U:H2'	23:BB:1507:C:C6	2.45	0.52
23:BB:1973:G:H2'	23:BB:1974:C:H6	1.75	0.52
23:BB:2014:A:OP1	40:BS:95:ARG:HB2	2.09	0.52
23:BB:2137:U:H6	23:BB:2137:U:O5'	1.93	0.52
23:BB:2298:A:H2'	23:BB:2299:U:O4'	2.10	0.52
23:BB:464:U:H2'	23:BB:465:G:O4'	2.10	0.52
23:BB:576:U:H2'	23:BB:577:G:C8	2.45	0.52
25:BC:112:GLY:O	25:BC:113:ASP:HB2	2.08	0.52
25:BC:114:GLN:HB2	25:BC:124:LYS:HZ2	1.73	0.52
25:BC:154:ALA:C	25:BC:156:SER:H	2.13	0.52
25:BC:20:ASN:HB2	25:BC:203:VAL:CG1	2.39	0.52
23:BB:1658:C:OP1	26:BD:136:ASN:HA	2.10	0.52
26:BD:165:MET:HG3	26:BD:166:GLY:H	1.74	0.52
27:BE:141:MET:HG3	27:BE:143:LEU:HB2	1.91	0.52
28:BF:18:GLU:C	28:BF:20:ASN:H	2.12	0.52
33:BL:116:VAL:HG11	33:BL:134:ALA:O	2.08	0.52
33:BL:63:LYS:O	33:BL:63:LYS:HD3	2.10	0.52
34:BM:33:LEU:HD13	34:BM:124:LEU:HD11	1.92	0.52
37:BP:93:LYS:HB2	37:BP:96:LEU:HA	1.91	0.52
39:BR:77:PHE:C	39:BR:78:ARG:HG3	2.30	0.52
42:BU:60:LYS:HG3	42:BU:61:GLU:N	2.24	0.52
42:BU:11:ILE:HB	42:BU:69:VAL:CG1	2.40	0.52
24:BV:42:LEU:HD12	24:BV:47:VAL:HG11	1.90	0.52
43:BW:66:VAL:HG13	43:BW:67:LYS:HG3	1.90	0.52
1:CA:1009:U:H2'	1:CA:1010:U:C6	2.44	0.52
1:CA:1070:U:H2'	1:CA:1071:C:H6	1.74	0.52
1:CA:1137:C:H1'	1:CA:1138:G:C2	2.44	0.52
1:CA:777:A:H2'	1:CA:778:G:C8	2.43	0.52
7:CH:111:THR:H	7:CH:114:ALA:HB3	1.74	0.52
11:CL:8:ARG:O	11:CL:10:PRO:HD3	2.08	0.52
11:CL:6:LEU:HB3	16:CQ:33:TYR:OH	2.09	0.52
22:DA:52:A:H2'	22:DA:53:A:H8	1.74	0.52
23:DB:110:G:O2'	23:DB:111:A:H5'	2.09	0.52
23:DB:1857:G:H2'	23:DB:1884:G:N2	2.24	0.52
23:DB:2364:C:H2'	23:DB:2365:G:O4'	2.10	0.52
23:DB:2882:A:OP1	35:DN:96:ARG:HD2	2.10	0.52
25:DC:172:THR:O	25:DC:173:LEU:CB	2.56	0.52
25:DC:193:GLU:C	25:DC:194:VAL:HG22	2.30	0.52
26:DD:122:VAL:HA	26:DD:128:ARG:CD	2.39	0.52
28:DF:55:ASP:O	28:DF:59:ILE:HB	2.10	0.52
23:DB:1099:G:O5'	52:DI:3:LYS:CA	2.57	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DI:73:PRO:CG	52:DI:78:LEU:HD21	2.39	0.52
32:DK:15:GLY:HA3	32:DK:52:VAL:HB	1.92	0.52
33:DL:59:ARG:HH21	50:D3:5:THR:HB	1.75	0.52
33:DL:89:VAL:HG21	33:DL:123:ARG:CZ	2.39	0.52
35:DN:123:GLU:CG	35:DN:124:ALA:H	2.23	0.52
36:DO:27:VAL:HG13	36:DO:40:ILE:HD11	1.91	0.52
37:DP:25:VAL:O	37:DP:47:ILE:HB	2.10	0.52
38:DQ:82:LEU:CD1	38:DQ:112:ALA:HB2	2.39	0.52
39:DR:78:ARG:HD2	39:DR:88:GLY:O	2.09	0.52
40:DS:28:LYS:HB3	40:DS:31:GLN:HB2	1.92	0.52
23:DB:1341:G:H1'	41:DT:59:ASN:HB3	1.90	0.52
42:DU:82:VAL:HG21	42:DU:95:PHE:O	2.09	0.52
43:DW:38:ARG:HE	43:DW:40:ARG:HA	1.75	0.52
1:AA:1414:U:H2'	1:AA:1415:G:H8	1.74	0.52
1:AA:663:A:O2'	1:AA:664:G:H5'	2.09	0.52
1:AA:78:A:O5'	1:AA:78:A:H8	1.93	0.52
1:AA:986:U:H1'	18:AS:53:GLY:O	2.10	0.52
3:AD:159:GLU:HG3	3:AD:160:LEU:HD13	1.91	0.52
6:AG:62:GLU:O	6:AG:66:GLU:HG3	2.10	0.52
9:AJ:9:ARG:C	9:AJ:10:LEU:HD12	2.30	0.52
13:AN:65:GLN:HE21	13:AN:65:GLN:HA	1.74	0.52
49:B2:8:SER:OG	49:B2:11:LYS:HD3	2.09	0.52
23:BB:1374:G:H2'	23:BB:1375:U:C6	2.45	0.52
23:BB:1534:U:O2'	23:BB:1535:A:H8	1.93	0.52
23:BB:1656:C:H2'	23:BB:1657:U:H6	1.74	0.52
23:BB:170:U:O2'	23:BB:171:U:H5'	2.09	0.52
23:BB:2405:G:H1'	23:BB:2412:A:H61	1.75	0.52
23:BB:2742:G:OP1	51:B4:36:ARG:NE	2.43	0.52
23:BB:540:C:O2'	23:BB:541:A:H5'	2.10	0.52
23:BB:554:U:H2'	23:BB:555:G:O4'	2.10	0.52
23:BB:580:U:H2'	23:BB:581:C:H6	1.70	0.52
23:BB:65:U:H2'	23:BB:66:C:C6	2.44	0.52
23:BB:991:C:H2'	23:BB:992:C:H6	1.74	0.52
25:BC:124:LYS:HB2	25:BC:125:PRO:HD2	1.91	0.52
25:BC:24:HIS:CB	25:BC:27:LYS:HE3	2.39	0.52
26:BD:101:PHE:O	26:BD:102:ALA:HB3	2.10	0.52
26:BD:4:LEU:H	26:BD:49:GLN:NE2	2.08	0.52
33:BL:29:LYS:CD	33:BL:30:THR:H	2.18	0.52
34:BM:1:MET:HA	34:BM:65:ILE:HG12	1.90	0.52
36:BO:19:GLN:HA	43:BW:76:ARG:NH2	2.25	0.52
36:BO:30:ARG:HG2	36:BO:30:ARG:HH11	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BO:55:GLU:HB2	36:BO:60:GLU:O	2.10	0.52
38:BQ:57:ARG:HA	38:BQ:60:TRP:CE3	2.43	0.52
39:BR:43:ASN:C	39:BR:45:GLU:H	2.11	0.52
39:BR:78:ARG:CB	39:BR:87:GLN:HA	2.31	0.52
24:BV:21:ARG:HE	24:BV:87:GLN:CB	2.23	0.52
24:BV:80:HIS:HB3	24:BV:83:LYS:O	2.09	0.52
43:BW:22:VAL:HG12	43:BW:23:LYS:N	2.25	0.52
46:BZ:64:PHE:CA	46:BZ:67:PRO:HD2	2.39	0.52
1:CA:1074:G:H2'	1:CA:1075:U:H6	1.74	0.52
1:CA:1149:C:O2'	1:CA:1150:A:H5'	2.09	0.52
1:CA:1307:U:H2'	1:CA:1308:U:C6	2.45	0.52
1:CA:1351:U:O2'	1:CA:1352:C:H5'	2.10	0.52
1:CA:33:A:H2'	1:CA:34:C:C6	2.45	0.52
1:CA:474:G:H2'	1:CA:475:C:H6	1.73	0.52
1:CA:711:G:O2'	1:CA:712:A:H5'	2.10	0.52
1:CA:730:G:O2'	1:CA:766:A:H5'	2.08	0.52
1:CA:82:G:H2'	1:CA:83:C:O4'	2.09	0.52
2:CC:149:LYS:HG3	2:CC:166:TRP:NE1	2.24	0.52
3:CD:18:LEU:HD12	3:CD:63:ILE:HB	1.91	0.52
5:CF:92:THR:O	5:CF:93:LYS:HB2	2.09	0.52
1:CA:939:G:H5''	6:CG:101:ARG:CZ	2.40	0.52
9:CJ:55:PRO:HA	13:CN:80:ARG:NH2	2.24	0.52
14:CO:62:ARG:NH1	14:CO:86:LEU:HD21	2.25	0.52
51:D4:11:CYS:SG	51:D4:26:ILE:C	2.88	0.52
51:D4:18:LYS:HE2	51:D4:19:ARG:H	1.68	0.52
23:DB:1050:A:H2'	23:DB:1051:G:O4'	2.09	0.52
23:DB:1133:A:H2	23:DB:2038:G:H21	1.58	0.52
23:DB:1309:G:H5'	49:D2:7:PRO:HB2	1.90	0.52
23:DB:137:U:C2'	23:DB:138:U:H5'	2.39	0.52
23:DB:1506:U:H2'	23:DB:1507:C:C6	2.45	0.52
23:DB:1739:A:H2'	23:DB:1740:G:O4'	2.10	0.52
23:DB:1986:C:O2'	23:DB:1987:A:H5'	2.10	0.52
23:DB:2083:G:H2'	23:DB:2084:C:H6	1.75	0.52
23:DB:2719:G:H4'	23:DB:2846:G:O3'	2.09	0.52
23:DB:448:U:O4'	23:DB:448:U:O2	2.27	0.52
23:DB:589:U:H2'	23:DB:590:A:H8	1.72	0.52
23:DB:737:C:O2'	23:DB:738:G:H5'	2.09	0.52
23:DB:931:U:H3	23:DB:1166:G:N2	2.07	0.52
26:DD:139:SER:O	26:DD:141:ARG:N	2.42	0.52
26:DD:31:ALA:HA	26:DD:51:THR:CA	2.34	0.52
27:DE:46:GLN:NE2	27:DE:87:ALA:N	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DG:41:GLU:HB2	29:DG:52:GLY:HA3	1.92	0.52
31:DJ:10:THR:O	31:DJ:11:VAL:HB	2.08	0.52
31:DJ:120:ARG:HB3	31:DJ:121:LYS:HZ1	1.74	0.52
31:DJ:49:ASP:O	31:DJ:50:THR:HB	2.09	0.52
33:DL:19:LEU:N	33:DL:19:LEU:HD22	2.24	0.52
23:DB:1252:G:H21	38:DQ:32:ARG:NE	2.07	0.52
38:DQ:47:ARG:HH22	38:DQ:50:ARG:HB2	1.75	0.52
39:DR:4:VAL:HG11	39:DR:46:GLU:OE2	2.10	0.52
39:DR:6:GLN:CG	39:DR:7:SER:H	2.18	0.52
40:DS:7:HIS:CD2	40:DS:46:LEU:HD13	2.44	0.52
24:DV:44:HIS:C	24:DV:46:LYS:H	2.13	0.52
43:DW:36:ILE:HD12	43:DW:37:VAL:H	1.73	0.52
43:DW:44:PHE:HB3	43:DW:77:LYS:O	2.09	0.52
36:DO:15:ARG:HH22	43:DW:76:ARG:CG	2.22	0.52
1:AA:1510:C:H2'	1:AA:1511:G:C8	2.44	0.52
1:AA:462:G:H2'	1:AA:463:U:C6	2.44	0.52
1:AA:764:C:C2'	1:AA:765:G:H5'	2.39	0.52
1:AA:894:G:O2'	1:AA:895:G:H5'	2.10	0.52
1:AA:924:C:O2'	1:AA:925:G:H5'	2.08	0.52
4:AE:35:LEU:HD11	4:AE:136:VAL:HG11	1.92	0.52
16:AQ:16:MET:CB	16:AQ:19:SER:HB2	2.39	0.52
21:AU:13:VAL:O	21:AU:14:ALA:HB2	2.09	0.52
51:B4:24:ARG:HG2	51:B4:26:ILE:H	1.74	0.52
23:BB:1149:G:H2'	23:BB:1150:C:H6	1.74	0.52
23:BB:1274:A:N3	23:BB:1297:C:H1'	2.25	0.52
23:BB:1341:G:H2'	23:BB:1397:U:O2'	2.09	0.52
23:BB:1717:A:H2'	23:BB:1718:G:O4'	2.10	0.52
23:BB:2070:A:H2'	23:BB:2071:A:H8	1.74	0.52
23:BB:107:G:H21	23:BB:346:A:H62	1.57	0.52
23:BB:445:C:H5'	27:BE:47:LYS:NZ	2.17	0.52
23:BB:27:G:H1'	23:BB:513:A:H61	1.74	0.52
23:BB:656:G:H2'	23:BB:657:U:C6	2.45	0.52
23:BB:920:A:H2'	23:BB:921:C:H6	1.75	0.52
25:BC:52:HIS:O	25:BC:53:ILE:HB	2.09	0.52
25:BC:61:TYR:CZ	25:BC:63:ILE:HD11	2.45	0.52
26:BD:154:LYS:C	26:BD:156:PHE:N	2.62	0.52
26:BD:34:VAL:HG23	26:BD:93:GLY:O	2.09	0.52
28:BF:103:ILE:HA	28:BF:107:VAL:HG21	1.90	0.52
52:BI:32:VAL:HG22	52:BI:60:VAL:CG2	2.40	0.52
32:BK:118:LEU:HD12	32:BK:118:LEU:N	2.24	0.52
32:BK:63:VAL:HG12	32:BK:103:VAL:HA	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BM:65:ILE:HG22	34:BM:101:VAL:HG12	1.92	0.52
40:BS:28:LYS:HB3	40:BS:71:VAL:CG2	2.40	0.52
24:BV:5:ASN:HA	24:BV:64:VAL:HB	1.92	0.52
43:BW:65:LYS:O	43:BW:66:VAL:HB	2.08	0.52
44:BX:47:ARG:O	44:BX:47:ARG:HG2	2.10	0.52
45:BY:3:THR:HA	45:BY:47:ILE:CD1	2.39	0.52
45:BY:9:THR:CB	45:BY:54:VAL:HA	2.34	0.52
1:CA:1213:A:O2'	1:CA:1214:C:H5''	2.10	0.52
1:CA:1317:C:H2'	1:CA:1318:A:O4'	2.10	0.52
1:CA:769:G:H4'	1:CA:1513:A:H4'	1.92	0.52
1:CA:303:A:H2'	1:CA:304:U:O4'	2.10	0.52
20:CB:104:LYS:NZ	20:CB:104:LYS:HB2	2.20	0.52
20:CB:44:LYS:O	20:CB:47:PRO:HD2	2.10	0.52
6:CG:49:LEU:HD23	6:CG:49:LEU:O	2.09	0.52
12:CM:100:ARG:HB2	12:CM:102:LYS:HG2	1.91	0.52
15:CP:22:ALA:CB	15:CP:32:PHE:HA	2.40	0.52
17:CR:23:LYS:C	17:CR:23:LYS:HD2	2.30	0.52
23:DB:686:U:H1'	49:D2:5:PHE:O	2.10	0.52
50:D3:46:LYS:HD2	50:D3:47:ALA:H	1.72	0.52
23:DB:1386:C:H2'	23:DB:1387:A:H8	1.74	0.52
23:DB:1745:A:O2'	23:DB:1746:A:H5'	2.09	0.52
23:DB:1845:G:O2'	23:DB:1846:G:H5'	2.10	0.52
23:DB:1957:C:H2'	23:DB:1958:C:H6	1.74	0.52
23:DB:2250:G:H21	23:DB:2496:C:C4'	2.22	0.52
23:DB:2316:G:O2'	23:DB:2317:A:H5'	2.10	0.52
23:DB:419:U:H2'	23:DB:420:C:H6	1.75	0.52
23:DB:600:G:H2'	23:DB:601:C:C6	2.44	0.52
23:DB:1818:U:H5''	25:DC:155:ARG:HG2	1.92	0.52
25:DC:231:HIS:ND1	25:DC:242:HIS:ND1	2.56	0.52
25:DC:179:GLU:OE2	25:DC:267:VAL:HG23	2.10	0.52
26:DD:129:THR:O	26:DD:140:HIS:HA	2.10	0.52
26:DD:4:LEU:CD1	26:DD:79:LEU:HD22	2.40	0.52
27:DE:147:LEU:HB2	27:DE:183:PHE:CD1	2.45	0.52
27:DE:21:ARG:HD3	27:DE:25:GLU:CB	2.40	0.52
55:DB:3322:HOH:O	27:DE:63:LYS:HE2	2.09	0.52
52:DI:105:LEU:HD11	52:DI:139:VAL:CG2	2.36	0.52
31:DJ:100:VAL:O	31:DJ:101:ILE:HB	2.10	0.52
33:DL:34:GLY:CA	39:DR:85:LYS:HE2	2.40	0.52
35:DN:14:SER:HA	35:DN:17:ARG:NH2	2.24	0.52
37:DP:52:ARG:O	37:DP:60:VAL:HG21	2.09	0.52
38:DQ:46:TYR:HA	39:DR:77:PHE:CD2	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DR:85:LYS:O	39:DR:86:GLN:HB3	2.10	0.52
41:DT:86:THR:O	41:DT:87:LEU:HB2	2.08	0.52
46:DZ:5:ILE:HG13	46:DZ:51:VAL:CG1	2.38	0.52
1:AA:1243:C:H2'	1:AA:1244:G:H8	1.75	0.52
1:AA:1518:A:H2'	1:AA:1519:A:C8	2.45	0.52
1:AA:217:C:O2'	1:AA:218:U:H5'	2.10	0.52
1:AA:541:G:H2'	1:AA:542:G:H8	1.75	0.52
1:AA:60:A:O2'	1:AA:61:G:P	2.66	0.52
1:AA:677:U:H2'	1:AA:678:U:C6	2.44	0.52
12:AM:53:ASP:HA	12:AM:56:ARG:CD	2.40	0.52
19:AT:57:VAL:C	19:AT:60:GLN:HE22	2.13	0.52
22:BA:32:U:H2'	22:BA:33:G:O4'	2.09	0.52
23:BB:52:A:C5	23:BB:118:A:C2	2.98	0.52
23:BB:1256:G:H21	27:BE:77:ILE:HG23	1.75	0.52
23:BB:1545:A:H2'	23:BB:1546:G:O4'	2.09	0.52
23:BB:1791:A:H5'	25:BC:207:ALA:HA	1.91	0.52
23:BB:1957:C:H2'	23:BB:1958:C:H6	1.74	0.52
23:BB:570:G:H2'	23:BB:2030:A:N7	2.24	0.52
23:BB:2189:U:H2'	23:BB:2190:G:H8	1.74	0.52
23:BB:2229:U:H2'	23:BB:2230:G:C8	2.45	0.52
23:BB:397:U:H2'	23:BB:398:C:C6	2.45	0.52
23:BB:807:U:O2'	23:BB:808:G:H5'	2.10	0.52
25:BC:141:HIS:HD2	25:BC:193:GLU:H	1.57	0.52
25:BC:230:PRO:O	25:BC:231:HIS:HB3	2.09	0.52
26:BD:115:GLY:HA3	26:BD:166:GLY:CA	2.39	0.52
52:BI:76:ALA:O	52:BI:80:LYS:HG3	2.10	0.52
31:BJ:75:TYR:N	31:BJ:75:TYR:CD2	2.78	0.52
37:BP:61:ARG:HB2	37:BP:61:ARG:CZ	2.40	0.52
41:BT:93:LEU:HD22	41:BT:93:LEU:N	2.23	0.52
42:BU:38:ILE:O	42:BU:39:ASN:HB2	2.08	0.52
1:CA:1226:C:H5''	12:CM:101:THR:O	2.10	0.52
1:CA:287:U:O2'	1:CA:288:A:H5'	2.08	0.52
1:CA:323:U:H2'	1:CA:324:G:O4'	2.10	0.52
4:CE:149:PRO:HG2	4:CE:150:GLU:OE2	2.09	0.52
9:CJ:9:ARG:CB	9:CJ:99:GLN:HG3	2.39	0.52
12:CM:68:LEU:HD22	12:CM:69:ARG:NH1	2.25	0.52
33:DL:62:PRO:HB3	50:D3:12:ARG:CD	2.40	0.52
50:D3:33:THR:HG22	50:D3:40:LYS:HD2	1.91	0.52
50:D3:40:LYS:O	50:D3:43:LEU:HD13	2.10	0.52
23:DB:1516:G:O2'	23:DB:1517:G:H5'	2.10	0.52
23:DB:2135:A:C2	23:DB:2136:G:C8	2.98	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2688:G:H1'	23:DB:2721:A:H61	1.74	0.52
23:DB:807:U:O2'	23:DB:808:G:H5'	2.10	0.52
31:DJ:41:LYS:CD	31:DJ:44:TYR:HB3	2.39	0.52
33:DL:123:ARG:CA	33:DL:142:ILE:HA	2.39	0.52
34:DM:41:LEU:CG	34:DM:46:ILE:HD11	2.40	0.52
35:DN:63:ARG:HA	35:DN:80:PHE:CZ	2.45	0.52
37:DP:50:ARG:HG3	37:DP:99:LEU:H	1.75	0.52
39:DR:44:GLY:H	39:DR:53:PHE:HE2	1.57	0.52
39:DR:76:LYS:HB3	39:DR:90:ARG:CB	2.40	0.52
41:DT:38:ALA:O	41:DT:42:GLU:HB3	2.10	0.52
41:DT:65:GLY:HA3	41:DT:76:ARG:NH2	2.25	0.52
43:DW:60:ALA:CB	43:DW:80:SER:HA	2.37	0.52
44:DX:6:LEU:HD22	44:DX:56:LEU:HD12	1.92	0.52
46:DZ:36:VAL:O	46:DZ:36:VAL:HG23	2.10	0.52
1:AA:1410:A:H2'	1:AA:1411:C:C6	2.45	0.52
1:AA:442:G:H2'	1:AA:443:C:C6	2.45	0.52
2:AC:154:GLY:HA2	2:AC:163:ARG:H	1.73	0.52
2:AC:30:ASP:O	2:AC:33:ASP:HB3	2.09	0.52
2:AC:61:LYS:O	2:AC:96:VAL:HB	2.09	0.52
2:AC:87:ARG:HG3	2:AC:100:ILE:CG2	2.38	0.52
3:AD:99:ASN:ND2	3:AD:110:ARG:HE	2.08	0.52
6:AG:41:ILE:HG13	6:AG:116:ALA:HA	1.91	0.52
7:AH:12:ARG:NH1	7:AH:26:MET:HB2	2.24	0.52
17:AR:70:THR:HG23	17:AR:71:ASP:H	1.75	0.52
23:BB:1213:A:C6	23:BB:1237:A:H1'	2.45	0.52
23:BB:1845:G:O2'	23:BB:1846:G:H5'	2.10	0.52
23:BB:2077:A:H2'	23:BB:2078:C:H6	1.72	0.52
23:BB:2186:G:H2'	23:BB:2187:U:O4'	2.10	0.52
23:BB:2354:C:O3'	43:BW:30:VAL:HG13	2.10	0.52
23:BB:2373:G:H2'	23:BB:2374:C:C6	2.45	0.52
23:BB:396:G:H5'	46:BZ:9:TYR:CB	2.34	0.52
23:BB:416:U:H2'	23:BB:417:C:H6	1.75	0.52
23:BB:526:A:N6	23:BB:2626:C:C4'	2.73	0.52
23:BB:753:A:H2'	23:BB:754:U:C6	2.45	0.52
23:BB:773:U:H5'	23:BB:774:G:OP2	2.10	0.52
23:BB:973:A:OP1	23:BB:973:A:H8	1.92	0.52
23:BB:978:G:H2'	23:BB:979:A:H8	1.75	0.52
25:BC:140:VAL:HG11	25:BC:163:ILE:HD11	1.92	0.52
25:BC:164:VAL:HB	25:BC:167:ASP:CG	2.30	0.52
25:BC:51:ARG:CD	25:BC:51:ARG:O	2.58	0.52
28:BF:136:ILE:HD12	28:BF:136:ILE:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:135:ILE:HB	28:BF:142:TYR:CE1	2.45	0.52
31:BJ:10:THR:HG22	31:BJ:50:THR:HA	1.90	0.52
23:BB:558:U:P	31:BJ:113:PRO:HG2	2.50	0.52
31:BJ:1:MET:SD	39:BR:14:VAL:HG21	2.50	0.52
34:BM:67:VAL:HB	34:BM:100:LYS:CD	2.40	0.52
35:BN:122:ALA:HA	47:B0:49:ARG:NH1	2.25	0.52
37:BP:63:ILE:N	37:BP:63:ILE:HD12	2.24	0.52
40:BS:11:ARG:HA	40:BS:100:THR:HG22	1.92	0.52
24:BV:82:TYR:HD2	24:BV:82:TYR:N	2.08	0.52
1:CA:1354:U:H2'	1:CA:1355:G:C8	2.44	0.52
1:CA:1355:G:O2'	1:CA:1356:G:H5'	2.10	0.52
1:CA:394:G:H2'	1:CA:395:C:C6	2.45	0.52
1:CA:537:G:H2'	1:CA:538:G:C8	2.45	0.52
20:CB:161:PHE:HA	20:CB:183:PHE:O	2.09	0.52
2:CC:156:LEU:HD12	2:CC:163:ARG:HG3	1.91	0.52
6:CG:96:ASN:O	6:CG:100:MET:HG2	2.10	0.52
8:CI:40:ARG:HA	8:CI:44:ARG:CZ	2.40	0.52
8:CI:9:GLY:HA2	8:CI:80:HIS:HB3	1.92	0.52
10:CK:85:VAL:CG2	10:CK:92:ARG:HH12	2.22	0.52
13:CN:53:ASP:HA	13:CN:58:ARG:HH11	1.75	0.52
18:CS:11:ASP:HB2	18:CS:14:LEU:HD23	1.92	0.52
19:CT:74:HIS:O	19:CT:78:LEU:HG	2.10	0.52
21:CU:16:ARG:HA	21:CU:19:LYS:NZ	2.25	0.52
49:D2:7:PRO:O	49:D2:8:SER:HB3	2.09	0.52
23:DB:1010:A:H4'	38:DQ:75:TYR:CD2	2.45	0.52
23:DB:1013:C:H2'	23:DB:1014:A:C8	2.44	0.52
23:DB:1441:G:H2'	23:DB:1442:U:H6	1.73	0.52
23:DB:1595:C:O2'	23:DB:1596:A:H5'	2.10	0.52
23:DB:1599:U:H2'	23:DB:1600:C:H6	1.75	0.52
23:DB:1739:A:H2'	23:DB:1740:G:H8	1.74	0.52
23:DB:2547:A:H2'	23:DB:2548:U:C6	2.45	0.52
23:DB:2599:G:C8	25:DC:234:GLY:HA2	2.45	0.52
23:DB:2862:G:H2'	23:DB:2863:C:C6	2.45	0.52
23:DB:352:A:H3'	23:DB:353:C:C6	2.45	0.52
23:DB:1790:C:O2'	25:DC:207:ALA:HB2	2.09	0.52
33:DL:57:LEU:O	33:DL:58:TYR:HB2	2.10	0.52
36:DO:25:ARG:CB	36:DO:94:ARG:HH22	2.23	0.52
40:DS:85:ILE:C	40:DS:85:ILE:HD13	2.31	0.52
41:DT:19:LYS:NZ	41:DT:26:LYS:HE2	2.25	0.52
42:DU:13:LEU:HD11	42:DU:69:VAL:H	1.75	0.52
45:DY:52:PHE:HB3	45:DY:53:MET:CE	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DZ:1:MET:HA	46:DZ:9:TYR:CE1	2.44	0.52
1:AA:938:A:H1'	1:AA:1376:U:O2'	2.10	0.52
5:AF:47:LEU:HB3	17:AR:65:SER:OG	2.10	0.52
7:AH:84:ILE:N	7:AH:84:ILE:HD12	2.24	0.52
8:AI:15:ALA:O	8:AI:66:VAL:HG23	2.09	0.52
9:AJ:88:MET:C	9:AJ:90:LEU:H	2.13	0.52
12:AM:63:VAL:HG12	12:AM:68:LEU:HG	1.92	0.52
1:AA:1048:G:H4'	13:AN:2:LYS:HZ2	1.72	0.52
1:AA:625:U:H4'	15:AP:16:PHE:CE2	2.44	0.52
48:B1:22:THR:O	48:B1:23:THR:HG23	2.10	0.52
23:BB:1745:A:H2'	23:BB:1746:A:C8	2.45	0.52
23:BB:1857:G:H2'	23:BB:1884:G:N2	2.24	0.52
23:BB:1939:U:H6	23:BB:1939:U:H5'	1.75	0.52
23:BB:1986:C:O2'	23:BB:1987:A:H5'	2.10	0.52
23:BB:2181:U:H2'	23:BB:2182:U:O4'	2.10	0.52
23:BB:2581:G:OP2	23:BB:2581:G:N2	2.43	0.52
23:BB:2862:G:H2'	23:BB:2863:C:C6	2.45	0.52
23:BB:458:G:H1'	23:BB:459:U:H5	1.73	0.52
23:BB:781:A:OP1	25:BC:216:ARG:NH2	2.42	0.52
25:BC:28:PRO:O	25:BC:29:PHE:CB	2.58	0.52
27:BE:113:VAL:HA	27:BE:117:ARG:HD3	1.92	0.52
27:BE:133:LEU:HD12	27:BE:136:GLN:OE1	2.09	0.52
28:BF:133:GLU:OE1	28:BF:147:ARG:HD3	2.10	0.52
28:BF:109:ARG:HG2	28:BF:136:ILE:HG13	1.92	0.52
31:BJ:7:LYS:HZ3	31:BJ:48:VAL:CB	2.22	0.52
32:BK:40:LYS:HZ2	32:BK:89:ASN:HD21	1.58	0.52
33:BL:132:ARG:HH11	33:BL:132:ARG:HG2	1.75	0.52
33:BL:82:LEU:O	33:BL:83:ALA:HB3	2.09	0.52
36:BO:59:ALA:CB	36:BO:63:LYS:HB2	2.40	0.52
31:BJ:45:THR:HB	38:BQ:63:ARG:HH12	1.75	0.52
38:BQ:87:VAL:HB	39:BR:55:ASP:OD2	2.10	0.52
44:BX:22:LEU:HA	44:BX:47:ARG:HH12	1.75	0.52
1:CA:1060:U:H5''	9:CJ:53:ILE:HG12	1.91	0.52
1:CA:1073:U:O2'	1:CA:1074:G:H5'	2.10	0.52
1:CA:1405:G:O2'	1:CA:1406:U:H5'	2.10	0.52
1:CA:246:A:C2	1:CA:282:A:C5	2.98	0.52
1:CA:894:G:O2'	1:CA:895:G:H5'	2.10	0.52
1:CA:902:G:H2'	1:CA:903:G:H8	1.74	0.52
20:CB:57:ASN:HA	20:CB:60:ALA:HB3	1.90	0.52
2:CC:127:VAL:HB	2:CC:128:MET:HE3	1.91	0.52
2:CC:2:GLN:HE21	2:CC:2:GLN:CA	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:CE:15:ILE:HG22	4:CE:16:ALA:N	2.25	0.52
4:CE:89:THR:CG2	4:CE:90:GLY:N	2.73	0.52
8:CI:114:LYS:H	8:CI:114:LYS:HD2	1.75	0.52
8:CI:26:LYS:HG3	8:CI:61:ASP:OD2	2.10	0.52
8:CI:56:MET:HG3	8:CI:57:VAL:HG23	1.92	0.52
11:CL:64:SER:HB2	11:CL:81:ILE:HD12	1.92	0.52
15:CP:54:LEU:HD13	15:CP:80:LYS:NZ	2.25	0.52
19:CT:43:LYS:HB3	19:CT:86:ALA:CB	2.39	0.52
22:DA:109:A:H2'	22:DA:110:C:H6	1.75	0.52
22:DA:44:G:N3	22:DA:47:C:N4	2.57	0.52
23:DB:1309:G:H4'	49:D2:7:PRO:HB2	1.92	0.52
23:DB:2086:U:H2'	23:DB:2087:G:H8	1.72	0.52
23:DB:2090:A:O3'	46:DZ:49:ARG:HD3	2.10	0.52
23:DB:1801:A:H5'	23:DB:2203:U:O2'	2.09	0.52
23:DB:2370:G:H2'	23:DB:2371:G:O4'	2.10	0.52
23:DB:283:G:H2'	23:DB:284:U:H6	1.75	0.52
23:DB:31:C:O2'	23:DB:32:C:H5'	2.10	0.52
28:DF:32:LYS:HE2	28:DF:156:THR:OG1	2.09	0.52
29:DG:151:ARG:HB2	29:DG:151:ARG:NH1	2.21	0.52
31:DJ:41:LYS:HE2	31:DJ:46:PRO:HD3	1.91	0.52
32:DK:61:VAL:CG1	32:DK:87:LEU:HD11	2.39	0.52
32:DK:95:ILE:O	32:DK:95:ILE:HG12	2.10	0.52
35:DN:42:LYS:CA	35:DN:42:LYS:HE3	2.39	0.52
35:DN:86:ARG:CB	35:DN:86:ARG:HH11	2.20	0.52
37:DP:27:VAL:HG23	37:DP:86:LYS:CB	2.40	0.52
41:DT:76:ARG:CG	41:DT:77:ARG:N	2.70	0.52
43:DW:60:ALA:HB3	43:DW:80:SER:CA	2.39	0.52
43:DW:67:LYS:CD	43:DW:71:LYS:H	2.23	0.52
45:DY:1:ALA:HB1	45:DY:37:ARG:HB3	1.92	0.52
1:AA:1060:U:H4'	9:AJ:54:SER:HB2	1.92	0.51
1:AA:978:A:HO2'	1:AA:1322:C:H5	1.55	0.51
1:AA:140:U:H2'	1:AA:141:G:H8	1.73	0.51
1:AA:648:A:H2'	1:AA:649:A:C8	2.45	0.51
4:AE:104:ILE:CD1	4:AE:114:LEU:HB2	2.40	0.51
47:B0:35:GLU:CG	47:B0:36:LYS:N	2.73	0.51
50:B3:23:HIS:CD2	50:B3:47:ALA:HA	2.45	0.51
51:B4:15:LYS:HB2	51:B4:26:ILE:HD11	1.91	0.51
23:BB:1228:G:H2'	23:BB:1229:C:C6	2.45	0.51
23:BB:1430:G:H2'	23:BB:1431:A:H8	1.74	0.51
23:BB:189:G:H2'	23:BB:205:G:N2	2.26	0.51
23:BB:1916:A:H2'	23:BB:1917:U:C6	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2052:A:C5'	26:BD:148:GLN:H	2.21	0.51
23:BB:208:C:H2'	23:BB:209:C:C6	2.45	0.51
23:BB:2195:U:H2'	23:BB:2196:C:H6	1.76	0.51
23:BB:233:A:H61	23:BB:428:A:H61	1.57	0.51
23:BB:27:G:HO2'	23:BB:28:A:H8	1.56	0.51
23:BB:995:C:OP2	38:BQ:53:LYS:HE2	2.08	0.51
27:BE:132:LYS:HD3	27:BE:133:LEU:HD22	1.91	0.51
27:BE:152:GLU:H	27:BE:171:ASP:HB3	1.75	0.51
27:BE:47:LYS:N	27:BE:47:LYS:HD3	2.12	0.51
28:BF:77:LYS:NZ	28:BF:79:ARG:HE	2.08	0.51
30:BH:125:THR:OG1	30:BH:146:VAL:HB	2.10	0.51
31:BJ:122:LEU:HD12	31:BJ:123:LYS:N	2.24	0.51
31:BJ:81:ILE:HG23	31:BJ:82:GLY:N	2.25	0.51
32:BK:47:ILE:HG23	32:BK:48:PRO:CD	2.40	0.51
38:BQ:47:ARG:HH21	38:BQ:51:GLN:N	2.08	0.51
40:BS:2:GLU:OE1	40:BS:108:SER:HB3	2.10	0.51
43:BW:45:HIS:ND1	43:BW:64:GLY:HA3	2.24	0.51
46:BZ:5:ILE:N	46:BZ:5:ILE:HD12	2.17	0.51
1:CA:1023:U:H2'	1:CA:1024:G:H8	1.75	0.51
1:CA:252:U:H2'	1:CA:253:A:H8	1.74	0.51
1:CA:280:C:H1'	16:CQ:39:ARG:NH1	2.25	0.51
1:CA:940:C:H2'	1:CA:941:G:C8	2.45	0.51
1:CA:960:U:O2'	1:CA:961:U:P	2.68	0.51
20:CB:31:PHE:N	20:CB:41:ASN:HB2	2.25	0.51
3:CD:121:ALA:HA	3:CD:145:ARG:HG3	1.93	0.51
3:CD:2:ARG:O	3:CD:3:TYR:HB3	2.10	0.51
6:CG:63:VAL:CG1	6:CG:126:ALA:HB1	2.40	0.51
6:CG:4:ARG:HB2	6:CG:4:ARG:NH1	2.25	0.51
7:CH:45:ILE:HD13	7:CH:60:LEU:HD11	1.91	0.51
10:CK:70:ALA:HA	10:CK:73:VAL:HG22	1.91	0.51
11:CL:23:LEU:C	11:CL:25:ALA:H	2.13	0.51
13:CN:50:LEU:CD1	13:CN:51:PRO:HD3	2.40	0.51
13:CN:65:GLN:HG2	13:CN:82:LYS:HG3	1.92	0.51
23:DB:1021:A:H61	23:DB:1142:A:H61	1.52	0.51
23:DB:1563:U:H2'	23:DB:1564:C:C6	2.45	0.51
23:DB:1639:C:H2'	23:DB:1640:A:C5'	2.33	0.51
23:DB:192:C:C2'	23:DB:193:U:H5'	2.39	0.51
23:DB:2440:C:H5'	55:DB:3424:HOH:O	2.10	0.51
23:DB:2755:C:H2'	51:D4:19:ARG:NH2	2.25	0.51
23:DB:656:G:H2'	23:DB:657:U:C6	2.45	0.51
23:DB:771:G:O2'	23:DB:772:C:H5'	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:845:A:N1	23:DB:847:U:H1'	2.25	0.51
23:DB:86:G:OP1	42:DU:28:LEU:HD22	2.10	0.51
23:DB:2599:G:N7	25:DC:234:GLY:HA2	2.26	0.51
23:DB:1654:A:O2'	26:DD:118:PHE:O	2.27	0.51
26:DD:146:ILE:CG1	26:DD:155:VAL:HG13	2.40	0.51
26:DD:56:LYS:HZ2	26:DD:59:ARG:HB2	1.75	0.51
31:DJ:61:LYS:HA	31:DJ:61:LYS:HE3	1.92	0.51
32:DK:105:ARG:O	32:DK:108:ARG:HB3	2.10	0.51
33:DL:109:LYS:HG2	33:DL:126:ARG:HH11	1.74	0.51
38:DQ:38:VAL:HA	38:DQ:41:ALA:HB3	1.92	0.51
39:DR:97:LYS:O	39:DR:98:ILE:HB	2.10	0.51
35:DN:107:ASN:HD21	40:DS:40:ASN:HB3	1.75	0.51
40:DS:46:LEU:O	40:DS:50:VAL:HG22	2.10	0.51
1:AA:1284:C:H3'	1:AA:1285:A:C5'	2.35	0.51
1:AA:250:A:H1'	1:AA:252:U:C5	2.45	0.51
1:AA:35:G:H2'	1:AA:36:C:C6	2.46	0.51
1:AA:783:C:O2'	1:AA:784:A:H5'	2.10	0.51
3:AD:145:ARG:HB3	3:AD:147:LYS:HE3	1.92	0.51
1:AA:590:U:OP1	7:AH:30:LYS:HG2	2.10	0.51
8:AI:122:ARG:NH1	8:AI:122:ARG:HG3	2.21	0.51
11:AL:119:LYS:HG3	11:AL:119:LYS:O	2.11	0.51
11:AL:80:LEU:O	11:AL:97:VAL:HG23	2.09	0.51
13:AN:30:ILE:HD12	13:AN:30:ILE:N	2.19	0.51
16:AQ:75:VAL:HG23	16:AQ:76:ARG:N	2.25	0.51
17:AR:31:TYR:CG	17:AR:54:LEU:HD21	2.45	0.51
23:BB:1438:U:H5'	23:BB:1516:G:O2'	2.10	0.51
23:BB:1563:U:H2'	23:BB:1564:C:C6	2.45	0.51
23:BB:167:A:H2'	23:BB:168:G:O4'	2.10	0.51
25:BC:180:MET:HB2	25:BC:265:PHE:HB2	1.93	0.51
25:BC:250:GLN:O	25:BC:251:THR:HB	2.09	0.51
26:BD:18:ASP:OD2	26:BD:20:VAL:HG13	2.09	0.51
27:BE:173:THR:HA	27:BE:201:ALA:O	2.11	0.51
30:BH:37:VAL:H	30:BH:38:PRO:CD	2.11	0.51
52:BI:112:LYS:O	52:BI:116:MET:HG3	2.09	0.51
23:BB:1081:U:H5''	52:BI:126:ARG:HD2	1.92	0.51
37:BP:26:GLU:CA	37:BP:46:VAL:HG13	2.38	0.51
37:BP:28:LYS:O	37:BP:30:TRP:N	2.43	0.51
37:BP:52:ARG:CB	37:BP:60:VAL:HG11	2.25	0.51
38:BQ:75:TYR:CZ	38:BQ:79:ILE:HD11	2.44	0.51
40:BS:89:ALA:O	40:BS:90:LYS:HB2	2.10	0.51
41:BT:25:GLU:HG3	41:BT:26:LYS:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BT:34:VAL:CG1	41:BT:43:ILE:HD11	2.39	0.51
42:BU:5:ARG:CG	42:BU:6:ARG:N	2.73	0.51
1:CA:1165:U:H2'	1:CA:1166:G:O4'	2.10	0.51
7:CH:104:SER:O	7:CH:122:GLY:HA3	2.10	0.51
8:CI:22:PRO:HA	8:CI:60:LEU:HB3	1.92	0.51
15:CP:23:ASP:OD1	15:CP:25:ARG:HB2	2.10	0.51
17:CR:33:THR:HG21	17:CR:37:LYS:HG2	1.92	0.51
47:D0:29:VAL:CG2	47:D0:32:THR:H	2.24	0.51
23:DB:1107:G:O2'	23:DB:1108:U:H5'	2.10	0.51
23:DB:1161:C:H2'	23:DB:1162:G:H8	1.76	0.51
23:DB:1560:G:H2'	23:DB:1561:C:C6	2.45	0.51
23:DB:2066:C:O2'	23:DB:2067:G:H5'	2.11	0.51
23:DB:2336:A:C1'	23:DB:2337:G:OP1	2.56	0.51
23:DB:2877:G:O2'	23:DB:2878:U:H5'	2.10	0.51
23:DB:753:A:H2'	23:DB:754:U:C6	2.45	0.51
35:DN:14:SER:HA	35:DN:17:ARG:HH22	1.75	0.51
37:DP:100:ARG:H	37:DP:100:ARG:HD2	1.76	0.51
39:DR:76:LYS:HD2	39:DR:90:ARG:HB3	1.90	0.51
1:AA:1108:G:H5'	2:AC:175:HIS:CD2	2.46	0.51
1:AA:1464:U:H2'	1:AA:1465:A:C8	2.45	0.51
1:AA:1475:G:OP1	23:BB:1689:A:H1'	2.10	0.51
1:AA:1512:U:O2'	1:AA:1513:A:H5'	2.10	0.51
1:AA:531:U:H6	1:AA:531:U:H5'	1.75	0.51
2:AC:75:VAL:O	2:AC:75:VAL:HG12	2.10	0.51
6:AG:14:ASP:CG	6:AG:15:PRO:HD2	2.31	0.51
11:AL:22:ALA:HB2	11:AL:56:LEU:CD2	2.39	0.51
10:AK:111:ASP:O	17:AR:72:ARG:HD3	2.10	0.51
51:B4:13:ASN:HD22	51:B4:13:ASN:N	2.08	0.51
23:BB:1205:A:C4'	23:BB:1206:G:OP2	2.59	0.51
23:BB:1516:G:O2'	23:BB:1517:G:H5'	2.10	0.51
23:BB:1595:C:O2'	23:BB:1596:A:H5'	2.10	0.51
23:BB:163:C:O2	23:BB:163:C:H5'	2.10	0.51
23:BB:1812:U:H2'	23:BB:1813:G:H8	1.75	0.51
23:BB:2518:A:H2'	23:BB:2518:A:N3	2.25	0.51
23:BB:2869:G:H2'	23:BB:2870:C:H6	1.75	0.51
23:BB:869:G:H2'	23:BB:870:U:C6	2.45	0.51
25:BC:81:GLU:HB3	25:BC:90:ILE:HG22	1.93	0.51
26:BD:46:ARG:HA	26:BD:81:GLU:O	2.09	0.51
29:BG:53:PRO:HG2	29:BG:61:TRP:HA	1.92	0.51
52:BI:27:LEU:N	52:BI:27:LEU:HD23	2.18	0.51
33:BL:21:ARG:HB3	33:BL:21:ARG:HH11	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:587:C:N3	33:BL:38:GLN:CB	2.73	0.51
37:BP:64:SER:HA	37:BP:72:VAL:HG12	1.91	0.51
38:BQ:30:VAL:O	38:BQ:33:VAL:HG12	2.11	0.51
38:BQ:34:ALA:C	38:BQ:38:VAL:HG12	2.31	0.51
39:BR:51:VAL:HG12	39:BR:53:PHE:H	1.74	0.51
24:BV:23:ALA:O	24:BV:24:ASN:HB2	2.11	0.51
1:CA:1275:A:H2'	1:CA:1276:G:O4'	2.11	0.51
1:CA:1477:U:H2'	1:CA:1478:U:C6	2.46	0.51
1:CA:217:C:O2'	1:CA:218:U:H5'	2.10	0.51
1:CA:392:C:H2'	1:CA:393:A:H8	1.75	0.51
1:CA:442:G:H2'	1:CA:443:C:C6	2.45	0.51
1:CA:648:A:H2'	1:CA:649:A:C8	2.46	0.51
1:CA:73:C:H2'	1:CA:74:A:C8	2.45	0.51
1:CA:837:U:H2'	1:CA:838:G:H8	1.75	0.51
1:CA:864:A:H2'	1:CA:865:A:C8	2.46	0.51
1:CA:975:A:HO2'	1:CA:1358:U:H1'	1.74	0.51
2:CC:33:ASP:OD1	2:CC:37:LYS:HE2	2.10	0.51
7:CH:38:VAL:HG13	7:CH:39:LEU:HD22	1.93	0.51
10:CK:109:ILE:N	10:CK:109:ILE:HD12	2.25	0.51
11:CL:56:LEU:HD12	11:CL:60:PHE:HB2	1.92	0.51
47:D0:36:LYS:HB2	47:D0:41:HIS:ND1	2.25	0.51
47:D0:42:ILE:O	47:D0:42:ILE:HG23	2.11	0.51
23:DB:1050:A:O2'	23:DB:2752:C:H1'	2.10	0.51
23:DB:146:A:H2'	23:DB:147:C:C6	2.44	0.51
23:DB:1534:U:O2'	23:DB:1535:A:H8	1.93	0.51
23:DB:1745:A:H2'	23:DB:1746:A:C8	2.45	0.51
23:DB:1812:U:H2'	23:DB:1813:G:H8	1.75	0.51
23:DB:2018:G:H1'	38:DQ:32:ARG:NH2	2.24	0.51
23:DB:2033:A:H1'	23:DB:2035:G:OP2	2.11	0.51
23:DB:2065:C:H2'	23:DB:2066:C:H6	1.74	0.51
23:DB:2281:A:O2'	23:DB:2282:G:H5'	2.11	0.51
23:DB:2327:A:H2'	23:DB:2328:A:C8	2.46	0.51
23:DB:2484:G:H21	34:DM:118:LYS:HG2	1.75	0.51
28:DF:66:ILE:O	28:DF:66:ILE:HG22	2.11	0.51
29:DG:10:VAL:HG23	29:DG:47:ASN:O	2.10	0.51
30:DH:4:ILE:HD13	30:DH:4:ILE:N	2.25	0.51
52:DI:53:PRO:CG	52:DI:77:VAL:HG11	2.39	0.51
31:DJ:136:GLN:N	31:DJ:137:PRO:CD	2.73	0.51
31:DJ:98:GLU:CD	31:DJ:98:GLU:H	2.12	0.51
33:DL:109:LYS:HE2	33:DL:126:ARG:NH1	2.26	0.51
23:DB:2394:C:H5'	33:DL:64:PHE:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:61:GLY:HA3	34:DM:105:MET:HB2	1.91	0.51
34:DM:97:GLN:N	34:DM:98:PRO:HD3	2.25	0.51
36:DO:18:LEU:HD21	36:DO:43:ASN:HD22	1.74	0.51
37:DP:54:LEU:N	37:DP:54:LEU:HD12	2.25	0.51
37:DP:90:ALA:HB3	37:DP:111:GLU:HB2	1.91	0.51
40:DS:9:HIS:O	40:DS:10:ALA:HB3	2.09	0.51
41:DT:3:ARG:HB3	41:DT:5:GLU:CD	2.31	0.51
42:DU:41:VAL:O	42:DU:42:LYS:HB2	2.10	0.51
24:DV:30:ILE:HG12	24:DV:91:PHE:CB	2.38	0.51
43:DW:10:ARG:HA	43:DW:10:ARG:NH1	2.26	0.51
45:DY:10:ARG:HG2	45:DY:31:ILE:HG21	1.92	0.51
1:AA:1348:U:H4'	8:AI:121:ARG:HH11	1.76	0.51
1:AA:676:A:O2'	1:AA:677:U:H5'	2.10	0.51
1:AA:883:C:O2'	1:AA:884:U:H5'	2.09	0.51
20:AB:116:LEU:HD12	20:AB:143:LEU:HD12	1.93	0.51
20:AB:216:VAL:O	20:AB:220:VAL:HG23	2.11	0.51
3:AD:84:ASN:HD22	3:AD:87:GLU:N	2.01	0.51
4:AE:14:LEU:HA	4:AE:36:THR:HG22	1.91	0.51
6:AG:72:VAL:HB	6:AG:144:ALA:CB	2.41	0.51
11:AL:3:VAL:HG23	11:AL:4:ASN:OD1	2.10	0.51
21:AU:5:VAL:HG12	21:AU:6:ARG:N	2.25	0.51
23:BB:1330:C:H2'	23:BB:1331:G:H8	1.76	0.51
23:BB:1560:G:H2'	23:BB:1561:C:C6	2.45	0.51
23:BB:165:A:H2'	23:BB:166:U:O4'	2.11	0.51
23:BB:740:C:H5''	23:BB:1784:A:OP1	2.11	0.51
23:BB:2742:G:O2'	23:BB:2743:U:H5'	2.11	0.51
23:BB:2771:C:H2'	23:BB:2772:C:H6	1.76	0.51
23:BB:2784:U:H4'	26:BD:41:ALA:O	2.10	0.51
23:BB:642:U:H2'	23:BB:644:A:OP2	2.10	0.51
23:BB:786:C:O2'	23:BB:787:C:H5'	2.10	0.51
26:BD:16:THR:HG23	26:BD:17:GLU:N	2.13	0.51
26:BD:55:LYS:NZ	26:BD:55:LYS:HB2	2.25	0.51
26:BD:67:HIS:O	26:BD:71:ALA:HB3	2.10	0.51
29:BG:51:PHE:HB3	29:BG:64:ALA:HB1	1.91	0.51
31:BJ:57:LEU:HD12	31:BJ:128:ASN:C	2.31	0.51
33:BL:106:GLU:CD	33:BL:106:GLU:H	2.12	0.51
33:BL:111:ILE:H	33:BL:111:ILE:CD1	2.23	0.51
37:BP:23:ASP:O	37:BP:25:VAL:HG23	2.09	0.51
39:BR:18:GLN:HA	39:BR:99:THR:HA	1.92	0.51
43:BW:3:LYS:HD2	43:BW:3:LYS:N	2.26	0.51
46:BZ:3:LYS:HB2	46:BZ:51:VAL:CG2	2.29	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1006:G:H2'	1:CA:1007:U:C6	2.45	0.51
1:CA:1464:U:H2'	1:CA:1465:A:C8	2.45	0.51
1:CA:146:G:O2'	1:CA:147:G:H5'	2.11	0.51
1:CA:250:A:H1'	1:CA:252:U:C5	2.45	0.51
1:CA:251:G:H4'	1:CA:252:U:H5'	1.92	0.51
1:CA:333:U:H2'	1:CA:334:C:H6	1.74	0.51
1:CA:81:A:H2'	1:CA:82:G:C8	2.45	0.51
1:CA:987:G:H2'	1:CA:988:G:C8	2.45	0.51
3:CD:101:VAL:O	3:CD:104:MET:HB2	2.11	0.51
4:CE:80:LEU:CD2	4:CE:122:VAL:HG11	2.40	0.51
12:CM:79:LEU:CD2	12:CM:86:ARG:HE	2.23	0.51
14:CO:31:LEU:O	14:CO:35:ILE:HG12	2.10	0.51
15:CP:12:LYS:O	15:CP:13:LYS:HB2	2.10	0.51
1:CA:376:G:C5'	15:CP:5:ARG:HD3	2.40	0.51
19:CT:49:ALA:O	19:CT:52:GLU:HB3	2.10	0.51
23:DB:1150:C:O2'	23:DB:1151:A:H5'	2.11	0.51
23:DB:1360:G:H2'	23:DB:1361:G:C5'	2.41	0.51
23:DB:2053:G:O2'	23:DB:2054:A:H5'	2.11	0.51
23:DB:2688:G:H1'	23:DB:2721:A:N6	2.25	0.51
23:DB:396:G:H2'	23:DB:397:U:H6	1.75	0.51
23:DB:441:U:H2'	23:DB:442:G:C8	2.45	0.51
23:DB:919:U:H6	23:DB:919:U:O5'	1.94	0.51
27:DE:4:VAL:HA	27:DE:14:VAL:HG22	1.92	0.51
27:DE:77:ILE:HG23	27:DE:78:TRP:CD1	2.45	0.51
28:DF:135:ILE:HG23	28:DF:136:ILE:H	1.75	0.51
31:DJ:77:HIS:HB2	31:DJ:85:LYS:CE	2.39	0.51
35:DN:61:ALA:O	35:DN:64:ARG:HB3	2.09	0.51
24:DV:2:PHE:HB3	24:DV:50:MET:SD	2.50	0.51
1:AA:1460:C:H2'	1:AA:1461:G:O4'	2.11	0.51
1:AA:394:G:H2'	1:AA:395:C:H6	1.75	0.51
1:AA:499:A:C2	1:AA:546:A:N3	2.79	0.51
20:AB:132:GLU:O	20:AB:136:ARG:HG3	2.10	0.51
2:AC:150:VAL:HG12	2:AC:151:GLU:N	2.25	0.51
3:AD:112:GLU:OE2	3:AD:153:ARG:HD3	2.11	0.51
8:AI:30:ASN:ND2	8:AI:30:ASN:N	2.59	0.51
8:AI:34:LEU:HD11	8:AI:47:VAL:CG2	2.40	0.51
15:AP:20:VAL:HG21	15:AP:32:PHE:CD2	2.45	0.51
47:B0:12:ARG:C	47:B0:12:ARG:HD3	2.31	0.51
48:B1:34:GLU:O	48:B1:35:LEU:HB3	2.10	0.51
23:BB:161:A:H3'	23:BB:162:U:C5'	2.25	0.51
23:BB:2033:A:H1'	23:BB:2035:G:OP2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2093:G:N7	23:BB:2225:A:H2'	2.26	0.51
23:BB:2322:A:N6	23:BB:2333:A:H62	2.09	0.51
23:BB:2547:A:H2'	23:BB:2548:U:C6	2.44	0.51
23:BB:2718:G:O2'	23:BB:2719:G:H5'	2.09	0.51
25:BC:153:LEU:HD13	25:BC:161:VAL:HG23	1.93	0.51
23:BB:2073:C:H5''	25:BC:227:VAL:HG23	1.92	0.51
26:BD:167:ASN:N	26:BD:167:ASN:ND2	2.58	0.51
29:BG:137:LYS:HA	29:BG:140:ILE:CG1	2.40	0.51
31:BJ:56:VAL:HG12	31:BJ:57:LEU:N	2.25	0.51
31:BJ:58:ASN:O	31:BJ:126:ALA:HA	2.10	0.51
33:BL:61:LEU:O	50:B3:11:LYS:NZ	2.43	0.51
33:BL:82:LEU:HD21	33:BL:92:LEU:HD12	1.92	0.51
23:BB:2469:A:H5''	34:BM:55:ARG:HD2	1.93	0.51
34:BM:8:LYS:HB2	34:BM:12:MET:CE	2.41	0.51
22:BA:114:C:H4'	36:BO:49:VAL:HG11	1.91	0.51
38:BQ:99:VAL:HG23	38:BQ:100:PHE:CD1	2.44	0.51
43:BW:65:LYS:HE2	43:BW:65:LYS:C	2.31	0.51
43:BW:6:GLY:C	43:BW:8:SER:H	2.14	0.51
46:BZ:69:SER:HB3	46:BZ:70:LYS:HE3	1.92	0.51
1:CA:541:G:H2'	1:CA:542:G:H8	1.75	0.51
1:CA:602:A:O2'	1:CA:603:U:H5'	2.10	0.51
1:CA:629:A:H2'	1:CA:630:A:O4'	2.11	0.51
20:CB:69:VAL:HB	20:CB:162:VAL:HB	1.91	0.51
11:CL:23:LEU:CD2	11:CL:58:ASN:HD22	2.18	0.51
13:CN:15:LEU:O	13:CN:19:TYR:HB2	2.10	0.51
13:CN:51:PRO:CB	13:CN:54:SER:HB3	2.40	0.51
19:CT:38:ILE:HD13	19:CT:85:LEU:HD23	1.93	0.51
21:CU:14:ALA:HB3	21:CU:16:ARG:HH12	1.74	0.51
22:DA:25:U:OP1	22:DA:25:U:H3'	2.10	0.51
22:DA:29:A:OP1	22:DA:29:A:C8	2.64	0.51
23:DB:1029:A:H3'	23:DB:1030:C:H6	1.76	0.51
23:DB:1315:C:H2'	23:DB:1316:U:C6	2.46	0.51
23:DB:1656:C:H2'	23:DB:1657:U:H6	1.74	0.51
23:DB:2328:A:H2'	23:DB:2329:U:H6	1.71	0.51
23:DB:2700:A:H2'	23:DB:2701:U:H6	1.76	0.51
23:DB:394:C:O2'	23:DB:395:U:H5'	2.11	0.51
23:DB:720:U:H2'	23:DB:721:A:C8	2.46	0.51
25:DC:171:VAL:HG12	25:DC:183:VAL:C	2.31	0.51
25:DC:243:PRO:O	25:DC:244:VAL:HG23	2.11	0.51
27:DE:139:LYS:HA	27:DE:143:LEU:CD2	2.40	0.51
27:DE:47:LYS:O	27:DE:49:ARG:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:6:LYS:HZ3	27:DE:8:ALA:HB2	1.75	0.51
52:DI:105:LEU:CD1	52:DI:129:GLU:HG2	2.37	0.51
52:DI:72:THR:CG2	52:DI:112:LYS:HD2	2.40	0.51
31:DJ:53:TYR:HA	31:DJ:121:LYS:HB3	1.93	0.51
31:DJ:14:ASP:H	31:DJ:53:TYR:HD1	1.57	0.51
34:DM:127:LYS:HZ3	34:DM:129:THR:H	1.59	0.51
36:DO:106:LEU:HD23	36:DO:106:LEU:C	2.31	0.51
22:DA:114:C:O2'	36:DO:49:VAL:HG23	2.10	0.51
40:DS:3:THR:OG1	40:DS:107:VAL:HG23	2.10	0.51
41:DT:17:SER:H	41:DT:20:ALA:CB	2.22	0.51
41:DT:55:VAL:CG2	41:DT:56:GLU:H	2.15	0.51
42:DU:78:LYS:HD2	42:DU:96:LYS:CG	2.41	0.51
43:DW:44:PHE:HD2	43:DW:77:LYS:HB3	1.75	0.51
23:DB:61:C:OP2	44:DX:44:LYS:HB2	2.11	0.51
1:AA:1448:C:H2'	1:AA:1449:C:H6	1.74	0.51
1:AA:212:G:H2'	1:AA:213:G:C8	2.46	0.51
1:AA:537:G:H2'	1:AA:538:G:C8	2.45	0.51
20:AB:102:ASN:OD1	20:AB:105:THR:HG22	2.10	0.51
2:AC:41:TYR:HA	2:AC:44:LYS:HD3	1.92	0.51
3:AD:104:MET:SD	3:AD:179:GLY:HA3	2.51	0.51
4:AE:84:VAL:HG21	4:AE:142:GLY:O	2.11	0.51
6:AG:117:LEU:HD13	6:AG:117:LEU:O	2.11	0.51
7:AH:79:ARG:HB2	7:AH:80:PRO:HD2	1.92	0.51
18:AS:3:SER:O	18:AS:4:LEU:HG	2.11	0.51
33:BL:62:PRO:HB3	50:B3:29:ARG:NH2	2.25	0.51
23:BB:1230:A:H2'	23:BB:1231:U:C6	2.46	0.51
23:BB:1599:U:H2'	23:BB:1600:C:H6	1.75	0.51
23:BB:2262:U:H4'	23:BB:2328:A:C2	2.46	0.51
23:BB:2621:G:HO2'	26:BD:164:GLN:HB2	1.75	0.51
23:BB:811:U:O5'	33:BL:31:GLY:HA2	2.11	0.51
23:BB:999:U:O2'	23:BB:1000:A:H5'	2.11	0.51
23:BB:589:U:H1'	27:BE:88:ARG:NH2	2.26	0.51
28:BF:150:GLY:O	28:BF:151:LEU:HD12	2.11	0.51
30:BH:73:ASN:ND2	30:BH:74:ALA:N	2.56	0.51
30:BH:8:LYS:HB3	30:BH:14:SER:HA	1.93	0.51
52:BI:79:LEU:HD23	52:BI:108:ILE:CD1	2.41	0.51
32:BK:78:ARG:O	37:BP:71:ARG:HB3	2.10	0.51
34:BM:20:LEU:HB2	34:BM:38:ARG:HD3	1.93	0.51
55:BB:3366:HOH:O	35:BN:15:SER:HB3	2.11	0.51
35:BN:38:LEU:HB2	35:BN:109:PRO:CG	2.41	0.51
36:BO:53:THR:O	36:BO:62:LEU:HD23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:112:ARG:N	37:BP:112:ARG:HH11	2.08	0.51
37:BP:77:SER:HB2	37:BP:78:PRO:HD3	1.92	0.51
38:BQ:61:ILE:O	38:BQ:65:ASN:HB3	2.11	0.51
41:BT:30:ILE:HG22	41:BT:85:VAL:HG22	1.92	0.51
43:BW:44:PHE:HB3	43:BW:76:ARG:HA	1.92	0.51
43:BW:66:VAL:CG1	43:BW:67:LYS:H	2.15	0.51
46:BZ:5:ILE:HB	46:BZ:51:VAL:HG12	1.93	0.51
1:CA:1021:A:H2'	1:CA:1022:A:C1'	2.41	0.51
1:CA:1085:U:H3'	1:CA:1086:U:C5	2.45	0.51
1:CA:1325:C:O2'	1:CA:1326:U:H5'	2.10	0.51
1:CA:423:G:H2'	1:CA:424:G:O4'	2.11	0.51
1:CA:531:U:H5'	1:CA:531:U:H6	1.75	0.51
1:CA:697:U:H2'	1:CA:698:G:H5'	1.93	0.51
2:CC:85:LYS:HG3	2:CC:86:LEU:N	2.24	0.51
3:CD:197:HIS:O	3:CD:201:GLU:HG3	2.10	0.51
19:CT:68:LYS:HB2	19:CT:70:LYS:HD3	1.93	0.51
23:DB:2045:C:H5''	47:D0:14:MET:SD	2.50	0.51
49:D2:37:LYS:HD3	49:D2:37:LYS:C	2.31	0.51
23:DB:1973:G:H2'	23:DB:1974:C:H6	1.75	0.51
23:DB:2199:A:O2'	46:DZ:34:LEU:HD21	2.09	0.51
23:DB:2597:G:C5'	25:DC:239:PHE:HB2	2.39	0.51
23:DB:274:C:H2'	23:DB:275:C:H6	1.76	0.51
23:DB:2849:U:O4	37:DP:96:LEU:HD21	2.11	0.51
23:DB:576:U:H2'	23:DB:577:G:C8	2.45	0.51
23:DB:686:U:H5''	49:D2:11:LYS:HE3	1.92	0.51
26:DD:37:VAL:CB	26:DD:46:ARG:HB2	2.39	0.51
27:DE:137:LYS:HA	27:DE:137:LYS:HZ3	1.73	0.51
27:DE:147:LEU:C	27:DE:148:ILE:HG23	2.31	0.51
28:DF:13:LYS:CA	28:DF:16:MET:HB2	2.34	0.51
29:DG:171:LYS:HE2	29:DG:174:LYS:HG3	1.92	0.51
32:DK:2:ILE:HA	32:DK:33:ALA:H	1.75	0.51
33:DL:132:ARG:HH12	33:DL:140:GLY:HA2	1.74	0.51
23:DB:2466:C:O2	34:DM:118:LYS:HD3	2.09	0.51
44:DX:7:ARG:O	44:DX:7:ARG:HG3	2.11	0.51
1:AA:1272:G:H2'	1:AA:1273:C:C6	2.45	0.51
1:AA:522:C:H41	11:AL:49:ARG:NH2	2.09	0.51
1:AA:837:U:H2'	1:AA:838:G:H8	1.75	0.51
6:AG:64:ALA:HA	6:AG:127:ALA:HB2	1.92	0.51
7:AH:6:ILE:CG2	7:AH:76:ARG:HH12	2.24	0.51
10:AK:63:GLN:HB3	10:AK:94:SER:OG	2.11	0.51
14:AO:29:ALA:HA	14:AO:84:LEU:HD21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AP:67:ILE:CG1	15:AP:71:VAL:HG13	2.41	0.51
15:AP:73:ALA:O	15:AP:77:GLU:HG3	2.10	0.51
23:BB:1025:G:H8	23:BB:1025:G:OP1	1.94	0.51
23:BB:1535:A:H3'	23:BB:1536:C:C6	2.46	0.51
23:BB:1771:C:H2'	23:BB:1772:A:C8	2.45	0.51
23:BB:205:G:HO2'	23:BB:206:U:P	2.34	0.51
23:BB:2074:U:H2'	23:BB:2075:U:C6	2.46	0.51
23:BB:2329:U:H2'	23:BB:2330:G:H8	1.74	0.51
23:BB:2688:G:H1'	23:BB:2721:A:N6	2.25	0.51
23:BB:2886:A:C8	47:B0:27:LEU:HD22	2.46	0.51
23:BB:354:A:C3'	23:BB:355:U:H5''	2.40	0.51
25:BC:141:HIS:ND1	25:BC:142:ASN:N	2.59	0.51
25:BC:160:TYR:HA	25:BC:193:GLU:CG	2.41	0.51
26:BD:15:PHE:HD2	26:BD:18:ASP:HB2	1.75	0.51
29:BG:120:ILE:CG1	29:BG:140:ILE:HG22	2.40	0.51
52:BI:18:ASN:N	52:BI:19:PRO:CD	2.74	0.51
33:BL:77:ILE:HA	33:BL:101:ILE:HD13	1.93	0.51
33:BL:78:ARG:HH21	33:BL:82:LEU:HD23	1.74	0.51
34:BM:9:PHE:CG	34:BM:10:ARG:N	2.79	0.51
35:BN:29:VAL:HG13	35:BN:83:LEU:HD13	1.93	0.51
41:BT:50:LEU:N	41:BT:50:LEU:HD22	2.26	0.51
42:BU:42:LYS:O	42:BU:57:ILE:HG13	2.11	0.51
42:BU:94:PHE:CZ	42:BU:100:GLU:HA	2.46	0.51
24:BV:80:HIS:CD2	24:BV:83:LYS:HB2	2.45	0.51
46:BZ:30:HIS:H	46:BZ:48:GLN:CD	2.13	0.51
1:CA:1004:A:H3'	1:CA:1005:A:H8	1.75	0.51
1:CA:488:C:H2'	1:CA:489:C:H6	1.76	0.51
1:CA:767:A:H2'	1:CA:768:A:C8	2.46	0.51
20:CB:34:ARG:HE	20:CB:39:ILE:HG13	1.74	0.51
3:CD:94:GLU:HA	3:CD:99:ASN:ND2	2.26	0.51
8:CI:33:SER:HB2	8:CI:36:GLN:HG2	1.92	0.51
16:CQ:16:MET:HB3	16:CQ:19:SER:HB2	1.92	0.51
16:CQ:80:LYS:H	16:CQ:80:LYS:HD2	1.76	0.51
47:D0:41:HIS:CD2	47:D0:42:ILE:HG22	2.46	0.51
50:D3:21:PHE:H	50:D3:48:MET:HB2	1.75	0.51
51:D4:23:ILE:HD13	51:D4:24:ARG:H	1.74	0.51
23:DB:2373:G:H2'	23:DB:2374:C:C6	2.45	0.51
23:DB:718:A:H3'	23:DB:719:C:H6	1.76	0.51
23:DB:862:G:H2'	23:DB:863:A:O4'	2.11	0.51
23:DB:866:A:H61	23:DB:913:U:C1'	2.24	0.51
25:DC:257:ARG:HA	25:DC:261:ARG:NE	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:188:MET:SD	27:DE:190:ALA:HB2	2.51	0.51
27:DE:1:MET:HG3	27:DE:18:THR:HG1	1.76	0.51
27:DE:88:ARG:HB3	27:DE:89:PRO:CD	2.40	0.51
29:DG:172:GLU:OE1	51:D4:32:LYS:HB3	2.11	0.51
30:DH:144:VAL:HG12	30:DH:145:ASN:H	1.76	0.51
30:DH:147:VAL:HG12	30:DH:148:ALA:N	2.25	0.51
30:DH:35:LYS:HZ1	30:DH:37:VAL:HG23	1.76	0.51
31:DJ:81:ILE:HG12	31:DJ:83:GLY:H	1.76	0.51
31:DJ:98:GLU:HG3	31:DJ:126:ALA:CB	2.41	0.51
37:DP:59:THR:O	37:DP:60:VAL:HG13	2.10	0.51
38:DQ:30:VAL:HG12	38:DQ:31:TYR:H	1.74	0.51
38:DQ:52:ARG:HG2	38:DQ:53:LYS:N	2.26	0.51
39:DR:23:GLU:O	39:DR:96:VAL:HG21	2.11	0.51
42:DU:59:GLU:HG3	42:DU:62:ALA:CB	2.38	0.51
46:DZ:59:ARG:HD3	46:DZ:63:ARG:CD	2.41	0.51
1:AA:1422:G:O2'	1:AA:1423:G:H5'	2.11	0.51
1:AA:488:C:H2'	1:AA:489:C:H6	1.76	0.51
5:AF:47:LEU:HD13	5:AF:51:ILE:HG22	1.92	0.51
13:AN:10:VAL:HG12	13:AN:11:LYS:NZ	2.26	0.51
1:AA:132:C:H5''	19:AT:68:LYS:CD	2.40	0.51
48:B1:7:LYS:HD2	48:B1:8:ILE:O	2.11	0.51
49:B2:10:LEU:O	49:B2:10:LEU:HD23	2.11	0.51
50:B3:24:LYS:HA	50:B3:45:PRO:CB	2.35	0.51
23:BB:2756:U:OP2	51:B4:19:ARG:HD3	2.10	0.51
51:B4:22:VAL:HG21	51:B4:24:ARG:HH21	1.76	0.51
23:BB:10:A:O2'	23:BB:11:C:H5'	2.11	0.51
23:BB:1913:A:C1'	23:BB:1914:C:OP1	2.56	0.51
23:BB:2464:G:H2'	23:BB:2465:C:H6	1.75	0.51
23:BB:2849:U:H4'	23:BB:2868:A:C2	2.46	0.51
23:BB:2877:G:O2'	23:BB:2878:U:H5'	2.11	0.51
23:BB:350:G:H2'	23:BB:351:C:C6	2.45	0.51
23:BB:496:G:H2'	23:BB:497:A:O4'	2.10	0.51
25:BC:160:TYR:O	25:BC:161:VAL:HG23	2.11	0.51
23:BB:2680:U:H5'	26:BD:194:PRO:HA	1.92	0.51
27:BE:105:LEU:HD22	27:BE:177:PRO:CG	2.40	0.51
23:BB:2060:A:H62	27:BE:69:ARG:HH21	1.57	0.51
27:BE:84:THR:HG23	27:BE:85:PHE:N	2.26	0.51
28:BF:90:LEU:HD23	28:BF:94:ARG:HG2	1.93	0.51
31:BJ:100:VAL:O	31:BJ:102:GLU:N	2.44	0.51
31:BJ:13:ARG:HG3	31:BJ:13:ARG:NH1	2.22	0.51
35:BN:37:THR:HG1	35:BN:40:LYS:HB2	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BO:68:LYS:H	36:BO:68:LYS:CD	2.23	0.51
32:BK:76:VAL:O	37:BP:73:PHE:HA	2.11	0.51
38:BQ:92:LYS:O	38:BQ:94:LEU:HD13	2.11	0.51
40:BS:71:VAL:HG12	40:BS:71:VAL:O	2.11	0.51
42:BU:2:ALA:C	42:BU:3:LYS:HD3	2.31	0.51
42:BU:6:ARG:C	42:BU:8:ASP:H	2.12	0.51
46:BZ:15:SER:HB3	46:BZ:23:LYS:CE	2.41	0.51
1:CA:1007:U:H2'	1:CA:1008:U:H6	1.75	0.51
1:CA:1049:U:O2'	1:CA:1050:G:P	2.68	0.51
1:CA:1262:C:H2'	1:CA:1263:C:C6	2.46	0.51
1:CA:1306:A:H61	1:CA:1331:G:C1'	2.18	0.51
1:CA:1369:C:H2'	1:CA:1370:G:C8	2.46	0.51
1:CA:1391:U:H2'	1:CA:1392:G:H8	1.74	0.51
1:CA:212:G:H2'	1:CA:213:G:C8	2.46	0.51
2:CC:189:HIS:CD2	2:CC:194:VAL:HG22	2.45	0.51
14:CO:35:ILE:O	14:CO:39:GLN:HG2	2.11	0.51
15:CP:57:ILE:O	15:CP:61:VAL:HG23	2.11	0.51
1:CA:451:A:C5'	15:CP:70:ARG:HH22	2.17	0.51
18:CS:26:ASP:OD2	18:CS:46:LEU:HA	2.11	0.51
50:D3:2:LYS:CB	50:D3:2:LYS:HZ2	2.24	0.51
51:D4:25:VAL:O	51:D4:35:GLN:HB2	2.10	0.51
23:DB:1005:C:H2'	23:DB:1006:C:C6	2.46	0.51
23:DB:1535:A:H3'	23:DB:1536:C:C6	2.46	0.51
23:DB:1654:A:H4'	35:DN:1:MET:O	2.11	0.51
23:DB:1717:A:H2'	23:DB:1718:G:O4'	2.10	0.51
23:DB:1864:U:OP1	23:DB:2410:G:O2'	2.22	0.51
23:DB:2291:U:H2'	23:DB:2292:U:H6	1.71	0.51
23:DB:2393:U:H5'	33:DL:61:LEU:O	2.10	0.51
23:DB:2515:C:O2'	23:DB:2516:A:H5'	2.11	0.51
23:DB:327:G:O2'	23:DB:328:U:H5'	2.11	0.51
23:DB:630:G:H1	33:DL:69:ARG:NH1	2.09	0.51
23:DB:877:A:H61	23:DB:898:C:H2'	1.74	0.51
23:DB:956:G:N2	23:DB:959:A:H3'	2.25	0.51
23:DB:2820:A:C5	26:DD:197:THR:HB	2.45	0.51
26:DD:5:VAL:CG2	26:DD:28:GLU:HA	2.40	0.51
27:DE:146:VAL:HG11	27:DE:184:ASP:OD1	2.11	0.51
28:DF:163:GLU:HA	28:DF:166:ARG:CD	2.41	0.51
29:DG:148:ARG:HH21	29:DG:153:PRO:HD2	1.76	0.51
52:DI:79:LEU:HD12	52:DI:135:MET:SD	2.51	0.51
32:DK:112:PHE:O	32:DK:114:LYS:N	2.43	0.51
33:DL:14:LYS:HD2	33:DL:15:ALA:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:DL:19:LEU:N	33:DL:19:LEU:HD13	2.25	0.51
36:DO:86:GLY:O	36:DO:87:ILE:HD13	2.10	0.51
38:DQ:82:LEU:HD11	38:DQ:108:LEU:HD11	1.92	0.51
39:DR:33:VAL:H	39:DR:35:PHE:HD2	1.59	0.51
43:DW:20:LEU:HD11	43:DW:31:LEU:HB2	1.92	0.51
44:DX:28:LEU:HD13	44:DX:42:LEU:HD11	1.93	0.51
45:DY:15:ARG:HD2	45:DY:53:MET:SD	2.51	0.51
1:AA:394:G:H2'	1:AA:395:C:C6	2.45	0.51
1:AA:1101:A:N7	20:AB:170:ILE:HD12	2.25	0.51
2:AC:42:LEU:O	2:AC:46:LEU:HB2	2.11	0.51
17:AR:22:TYR:HB2	17:AR:61:ALA:HB2	1.92	0.51
47:B0:45:ASP:HB2	47:B0:47:TYR:CD1	2.46	0.51
49:B2:21:ARG:HB3	49:B2:31:LEU:HD21	1.91	0.51
51:B4:9:LYS:O	51:B4:25:VAL:HG13	2.11	0.51
23:BB:1315:C:H2'	23:BB:1316:U:C6	2.46	0.51
23:BB:1733:G:H2'	23:BB:1734:G:C8	2.45	0.51
23:BB:2364:C:H2'	23:BB:2365:G:O4'	2.10	0.51
23:BB:2666:C:O4'	23:BB:2666:C:O2	2.24	0.51
23:BB:1663:G:O2'	23:BB:2686:G:H4'	2.11	0.51
23:BB:320:A:H2'	27:BE:131:THR:HG21	1.92	0.51
23:BB:55:G:H2'	23:BB:56:A:H8	1.76	0.51
26:BD:29:VAL:HG23	26:BD:51:THR:HG21	1.93	0.51
27:BE:169:VAL:HG12	27:BE:171:ASP:OD1	2.11	0.51
52:BI:27:LEU:HB2	52:BI:32:VAL:HG21	1.92	0.51
38:BQ:103:VAL:O	38:BQ:106:THR:HG22	2.11	0.51
38:BQ:92:LYS:HA	38:BQ:95:ALA:CB	2.40	0.51
38:BQ:92:LYS:HD3	38:BQ:93:ILE:HG22	1.92	0.51
39:BR:40:MET:HA	39:BR:54:VAL:HG11	1.93	0.51
42:BU:69:VAL:HG13	42:BU:77:GLY:H	1.75	0.51
46:BZ:59:ARG:HH21	46:BZ:62:LYS:NZ	2.09	0.51
1:CA:1041:G:H2'	1:CA:1042:A:C8	2.44	0.51
1:CA:1238:A:C2	1:CA:1241:G:N3	2.79	0.51
1:CA:1287:A:O2'	1:CA:1288:A:H5'	2.11	0.51
1:CA:402:G:H2'	1:CA:403:C:H6	1.75	0.51
1:CA:538:G:P	11:CL:111:GLN:HB2	2.51	0.51
1:CA:499:A:C2	1:CA:546:A:N3	2.79	0.51
1:CA:769:G:O2'	1:CA:770:C:H5'	2.11	0.51
20:CB:132:GLU:HG2	20:CB:135:MET:HE1	1.93	0.51
3:CD:75:TYR:HE1	3:CD:200:VAL:HG23	1.75	0.51
3:CD:89:LEU:HA	3:CD:92:LEU:HB2	1.93	0.51
5:CF:53:LYS:HE2	5:CF:54:LEU:HD23	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:54:LEU:HD13	5:CF:55:HIS:H	1.76	0.51
5:CF:5:GLU:HA	5:CF:63:ASN:HA	1.92	0.51
8:CI:51:LEU:CB	8:CI:56:MET:HG2	2.38	0.51
12:CM:106:ARG:NH1	12:CM:109:LYS:HD2	2.26	0.51
23:DB:2017:U:H4'	47:D0:4:GLN:O	2.11	0.51
49:D2:16:HIS:NE2	49:D2:44:VAL:HG12	2.26	0.51
50:D3:12:ARG:HD2	50:D3:24:LYS:O	2.10	0.51
50:D3:4:LYS:CE	50:D3:61:LEU:HD13	2.41	0.51
23:DB:105:C:H2'	23:DB:106:C:H6	1.76	0.51
23:DB:1274:A:N3	23:DB:1297:C:H1'	2.25	0.51
23:DB:1367:A:H2'	23:DB:1368:G:H5'	1.93	0.51
23:DB:175:G:H2'	23:DB:176:A:H8	1.74	0.51
23:DB:2581:G:OP2	23:DB:2581:G:N2	2.43	0.51
23:DB:532:A:H5''	38:DQ:27:ARG:HH12	1.76	0.51
23:DB:633:A:O5'	23:DB:633:A:C8	2.57	0.51
23:DB:642:U:H2'	23:DB:644:A:OP2	2.10	0.51
26:DD:204:LYS:HB3	26:DD:205:PRO:CD	2.30	0.51
52:DI:23:VAL:HG12	52:DI:24:GLY:N	2.26	0.51
31:DJ:15:TRP:CE3	31:DJ:138:GLN:HB2	2.45	0.51
33:DL:41:ARG:HA	33:DL:41:ARG:CZ	2.41	0.51
34:DM:131:VAL:CG2	34:DM:133:LYS:HB3	2.40	0.51
23:DB:2293:G:P	36:DO:13:ARG:HH22	2.33	0.51
36:DO:35:ILE:HG21	36:DO:74:VAL:HG21	1.93	0.51
40:DS:68:ASP:HB2	40:DS:69:LEU:HD22	1.93	0.51
23:DB:2278:A:N6	43:DW:10:ARG:HB2	2.25	0.51
46:DZ:25:ARG:O	46:DZ:26:SER:HB3	2.11	0.51
46:DZ:50:ASP:HA	46:DZ:53:THR:HG22	1.92	0.51
1:AA:1122:U:H2'	1:AA:1123:U:C6	2.46	0.51
1:AA:1202:U:C2'	1:AA:1203:C:H5'	2.41	0.51
1:AA:1150:A:H1'	1:AA:1280:A:N6	2.26	0.51
1:AA:1338:G:H2'	1:AA:1339:A:C8	2.45	0.51
1:AA:1486:G:H2'	1:AA:1487:G:C8	2.46	0.51
1:AA:539:A:H2'	1:AA:540:G:H8	1.74	0.51
1:AA:767:A:H2'	1:AA:768:A:C8	2.46	0.51
20:AB:67:LEU:HD21	20:AB:157:PRO:CG	2.41	0.51
2:AC:146:LYS:HB2	2:AC:146:LYS:NZ	2.25	0.51
2:AC:184:ASN:HD22	2:AC:185:THR:N	2.08	0.51
8:AI:10:ARG:HA	8:AI:77:ALA:HB1	1.92	0.51
8:AI:113:LYS:HG2	8:AI:114:LYS:N	2.26	0.51
10:AK:110:THR:CG2	21:AU:4:LYS:HD2	2.39	0.51
23:BB:1269:A:H2'	23:BB:1270:C:C6	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1360:G:H2'	23:BB:1361:G:C5'	2.41	0.51
23:BB:2076:U:O4'	23:BB:2076:U:O2	2.29	0.51
23:BB:2520:C:C6	23:BB:2567:G:H1'	2.46	0.51
23:BB:2653:U:H3'	23:BB:2654:A:H2'	1.92	0.51
23:BB:538:A:O2'	23:BB:539:G:H5'	2.10	0.51
23:BB:854:C:O2'	23:BB:855:G:H5'	2.11	0.51
25:BC:248:GLY:HA3	25:BC:252:LYS:HE3	1.91	0.51
27:BE:119:ILE:HG23	27:BE:189:THR:O	2.11	0.51
23:BB:37:C:C2	27:BE:46:GLN:OE1	2.64	0.51
27:BE:78:TRP:HE3	27:BE:78:TRP:N	2.09	0.51
28:BF:108:PRO:HB3	28:BF:113:PHE:CD2	2.46	0.51
28:BF:121:PHE:CZ	28:BF:169:LEU:HD12	2.46	0.51
28:BF:33:ILE:HG13	28:BF:34:THR:H	1.76	0.51
23:BB:2314:A:H5'	28:BF:34:THR:HG21	1.92	0.51
36:BO:25:ARG:HD3	36:BO:40:ILE:HD11	1.92	0.51
37:BP:23:ASP:OD2	37:BP:93:LYS:HG2	2.10	0.51
41:BT:43:ILE:O	41:BT:47:VAL:HG23	2.11	0.51
41:BT:96:VAL:HG22	41:BT:97:GLY:H	1.75	0.51
1:CA:205:A:H2'	1:CA:206:C:C6	2.44	0.51
1:CA:35:G:H2'	1:CA:36:C:C6	2.46	0.51
1:CA:53:A:C2	1:CA:54:C:H1'	2.46	0.51
1:CA:662:U:H2'	1:CA:663:A:H8	1.75	0.51
4:CE:28:ARG:NH2	4:CE:30:PHE:HB3	2.26	0.51
4:CE:156:ARG:HH22	7:CH:100:ILE:HG23	1.75	0.51
10:CK:14:GLN:HA	10:CK:77:GLY:HA3	1.92	0.51
12:CM:50:GLY:O	12:CM:54:THR:HG23	2.11	0.51
12:CM:6:ILE:N	12:CM:6:ILE:HD12	2.26	0.51
23:DB:1196:C:H2'	23:DB:1197:G:H8	1.75	0.51
23:DB:1248:G:P	27:DE:44:ARG:HH22	2.34	0.51
23:DB:1309:G:OP1	49:D2:9:VAL:HG22	2.11	0.51
23:DB:2241:A:O2'	23:DB:2242:G:H5'	2.10	0.51
23:DB:2669:G:H2'	23:DB:2670:A:H8	1.76	0.51
23:DB:2821:A:H2'	23:DB:2822:G:C8	2.46	0.51
23:DB:361:G:O2'	23:DB:362:A:H5'	2.11	0.51
23:DB:545:U:H3'	23:DB:546:U:C5'	2.40	0.51
23:DB:919:U:H2'	23:DB:920:A:H8	1.69	0.51
23:DB:1828:G:O6	25:DC:219:VAL:HG11	2.11	0.51
25:DC:48:ILE:CG2	25:DC:49:THR:H	2.06	0.51
23:DB:779:U:P	25:DC:49:THR:HG1	2.34	0.51
26:DD:5:VAL:HG12	26:DD:6:GLY:H	1.75	0.51
27:DE:85:PHE:O	27:DE:86:ALA:CB	2.59	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DG:3:VAL:HG22	29:DG:4:ALA:N	2.25	0.51
30:DH:68:ARG:HD3	30:DH:71:LYS:HB2	1.92	0.51
52:DI:59:THR:O	52:DI:59:THR:HG23	2.09	0.51
31:DJ:123:LYS:N	31:DJ:123:LYS:HD2	2.25	0.51
33:DL:124:GLY:C	33:DL:125:LEU:HD23	2.32	0.51
33:DL:39:LYS:NZ	33:DL:39:LYS:H	2.08	0.51
34:DM:118:LYS:HB2	34:DM:118:LYS:HZ3	1.75	0.51
23:DB:910:A:H62	34:DM:15:GLY:HA3	1.75	0.51
24:DV:21:ARG:HH21	24:DV:88:HIS:N	2.08	0.51
1:AA:1498:U:H3'	53:AA:1601:KSG:HN12	1.77	0.50
1:AA:370:C:H2'	1:AA:371:A:C8	2.46	0.50
1:AA:692:U:H5	10:AK:27:ASN:HD22	1.58	0.50
1:AA:712:A:O2'	1:AA:713:G:H5'	2.12	0.50
1:AA:80:A:N3	1:AA:81:A:H1'	2.26	0.50
4:AE:55:VAL:N	4:AE:56:PRO:HD2	2.26	0.50
5:AF:7:VAL:HG11	17:AR:64:LEU:HD21	1.92	0.50
8:AI:38:PHE:C	8:AI:44:ARG:HG2	2.31	0.50
9:AJ:30:LYS:HD2	9:AJ:30:LYS:O	2.11	0.50
9:AJ:73:LEU:CD1	9:AJ:75:ASP:HB3	2.42	0.50
15:AP:36:VAL:HG13	15:AP:36:VAL:O	2.12	0.50
16:AQ:46:HIS:HB2	16:AQ:70:LYS:HE3	1.93	0.50
23:BB:1739:A:H2'	23:BB:1740:G:H8	1.74	0.50
23:BB:1797:G:H5'	25:BC:251:THR:HA	1.93	0.50
23:BB:2195:U:H2'	23:BB:2196:C:C6	2.46	0.50
25:BC:160:TYR:HE2	25:BC:194:VAL:HA	1.76	0.50
26:BD:129:THR:O	26:BD:130:GLN:HB2	2.10	0.50
26:BD:83:ARG:HG2	26:BD:83:ARG:HH11	1.76	0.50
27:BE:3:LEU:HA	27:BE:117:ARG:CZ	2.41	0.50
28:BF:118:ALA:HB2	28:BF:177:ARG:NH1	2.25	0.50
29:BG:63:GLN:HG3	29:BG:64:ALA:N	2.26	0.50
30:BH:57:LYS:HA	30:BH:57:LYS:HE3	1.93	0.50
32:BK:84:CYS:O	32:BK:85:VAL:HB	2.10	0.50
34:BM:73:ILE:O	34:BM:74:THR:HB	2.11	0.50
35:BN:120:GLU:C	35:BN:121:LYS:HD2	2.32	0.50
38:BQ:61:ILE:HB	38:BQ:75:TYR:CE2	2.46	0.50
39:BR:37:GLU:HG3	39:BR:38:VAL:N	2.26	0.50
41:BT:58:VAL:HG12	41:BT:59:ASN:N	2.26	0.50
24:BV:53:LYS:HE3	24:BV:55:GLU:HG3	1.92	0.50
43:BW:45:HIS:HA	43:BW:75:ASN:HB2	1.91	0.50
1:CA:1251:A:O2'	1:CA:1370:G:H5'	2.12	0.50
1:CA:1461:G:O2'	1:CA:1462:C:H5'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:176:C:H2'	1:CA:177:G:N3	2.26	0.50
1:CA:865:A:C2	1:CA:918:A:H4'	2.46	0.50
4:CE:104:ILE:C	4:CE:104:ILE:HD13	2.31	0.50
8:CI:29:ILE:O	8:CI:29:ILE:HG23	2.11	0.50
9:CJ:28:THR:HG22	9:CJ:86:ALA:HB1	1.92	0.50
1:CA:1228:C:C5'	12:CM:112:ARG:HB3	2.41	0.50
12:CM:23:GLY:N	12:CM:69:ARG:HH22	2.09	0.50
12:CM:85:TYR:HA	12:CM:88:LEU:HD12	1.92	0.50
14:CO:11:VAL:HG21	14:CO:21:THR:HG22	1.92	0.50
15:CP:60:TRP:HB3	15:CP:65:ALA:HB2	1.92	0.50
16:CQ:32:ILE:O	16:CQ:32:ILE:HG12	2.11	0.50
1:CA:1319:A:P	18:CS:4:LEU:HD21	2.51	0.50
18:CS:62:THR:HG23	18:CS:63:ASP:N	2.26	0.50
22:DA:40:U:O2	22:DA:43:C:H5''	2.11	0.50
23:DB:1083:U:H2'	23:DB:1085:A:OP2	2.11	0.50
23:DB:1091:G:O2'	23:DB:1092:C:H5'	2.11	0.50
23:DB:1175:A:C3'	23:DB:1176:U:H5'	2.40	0.50
23:DB:2313:C:H2'	23:DB:2314:A:H8	1.76	0.50
23:DB:2653:U:H3'	23:DB:2654:A:H2'	1.92	0.50
23:DB:466:A:H2'	23:DB:467:G:H5'	1.92	0.50
28:DF:36:ASN:O	28:DF:151:LEU:HD12	2.11	0.50
31:DJ:122:LEU:O	31:DJ:123:LYS:HB2	2.11	0.50
31:DJ:81:ILE:HG12	31:DJ:82:GLY:N	2.24	0.50
33:DL:103:ILE:HG13	33:DL:104:GLN:N	2.22	0.50
23:DB:2360:G:H5'	33:DL:58:TYR:OH	2.11	0.50
33:DL:59:ARG:HH12	50:D3:58:ILE:HG21	1.75	0.50
34:DM:15:GLY:C	34:DM:16:ARG:HG3	2.31	0.50
35:DN:102:PHE:CZ	35:DN:104:ALA:HB2	2.46	0.50
35:DN:9:GLN:HE22	35:DN:39:PRO:HB3	1.75	0.50
36:DO:90:VAL:HG11	36:DO:116:GLN:O	2.10	0.50
39:DR:4:VAL:HA	39:DR:43:ASN:CG	2.32	0.50
39:DR:69:GLY:H	39:DR:97:LYS:HB2	1.75	0.50
40:DS:45:VAL:C	40:DS:47:VAL:H	2.14	0.50
24:DV:26:PHE:CE1	24:DV:42:LEU:HD12	2.46	0.50
43:DW:42:THR:HG21	43:DW:66:VAL:HG13	1.93	0.50
46:DZ:32:LEU:N	46:DZ:32:LEU:HD22	2.26	0.50
46:DZ:55:GLY:O	46:DZ:60:PHE:N	2.44	0.50
46:DZ:64:PHE:C	46:DZ:67:PRO:HD2	2.31	0.50
1:AA:1402:C:H2'	1:AA:1403:C:O4'	2.11	0.50
1:AA:402:G:H2'	1:AA:403:C:H6	1.75	0.50
1:AA:505:G:H5'	1:AA:534:U:H2'	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:53:A:C2	1:AA:54:C:H1'	2.46	0.50
1:AA:629:A:H2'	1:AA:630:A:O4'	2.11	0.50
20:AB:139:GLU:HG2	20:AB:143:LEU:HG	1.92	0.50
2:AC:38:VAL:HG12	2:AC:93:ILE:CG2	2.42	0.50
3:AD:170:LEU:HB2	3:AD:180:THR:O	2.11	0.50
8:AI:32:ARG:NH2	8:AI:36:GLN:HG2	2.26	0.50
10:AK:113:THR:HG21	21:AU:28:LEU:HD21	1.93	0.50
21:AU:3:ILE:HG12	21:AU:19:LYS:HB3	1.92	0.50
50:B3:47:ALA:O	50:B3:48:MET:HG3	2.11	0.50
23:BB:243:U:OP1	50:B3:4:LYS:HG2	2.11	0.50
23:BB:1439:A:N7	23:BB:1440:U:C2	2.79	0.50
23:BB:2515:C:O2'	23:BB:2516:A:H5'	2.11	0.50
23:BB:256:A:H2'	23:BB:257:C:H6	1.77	0.50
23:BB:2636:C:H2'	23:BB:2637:U:C6	2.46	0.50
23:BB:221:A:H2'	23:BB:266:G:N7	2.26	0.50
23:BB:2896:C:H2'	23:BB:2897:U:C6	2.46	0.50
23:BB:557:C:H2'	23:BB:558:U:C6	2.46	0.50
23:BB:841:G:O2'	23:BB:842:U:H5'	2.12	0.50
23:BB:85:G:OP1	42:BU:5:ARG:HA	2.12	0.50
25:BC:172:THR:O	25:BC:173:LEU:HB2	2.11	0.50
25:BC:207:ALA:HA	25:BC:211:ARG:HB2	1.93	0.50
25:BC:266:ILE:HG13	25:BC:266:ILE:O	2.10	0.50
25:BC:44:ASN:HA	25:BC:49:THR:O	2.11	0.50
26:BD:13:ARG:HB2	26:BD:21:SER:HB2	1.93	0.50
26:BD:17:GLU:OE1	37:BP:32:VAL:HG11	2.12	0.50
26:BD:87:GLY:O	26:BD:88:GLU:HB2	2.10	0.50
27:BE:151:GLY:HA2	27:BE:171:ASP:HA	1.93	0.50
28:BF:140:ILE:HG22	28:BF:142:TYR:H	1.76	0.50
28:BF:176:PHE:O	28:BF:177:ARG:HB3	2.11	0.50
31:BJ:137:PRO:O	31:BJ:139:VAL:HG13	2.11	0.50
33:BL:126:ARG:NH2	33:BL:131:ALA:CB	2.73	0.50
38:BQ:22:GLY:HA2	38:BQ:28:SER:OG	2.10	0.50
38:BQ:71:ASN:HD22	38:BQ:109:VAL:HG21	1.77	0.50
38:BQ:84:LYS:O	38:BQ:85:ALA:HB2	2.12	0.50
39:BR:4:VAL:HG22	39:BR:5:PHE:N	2.26	0.50
42:BU:73:ASN:HB2	42:BU:96:LYS:HZ2	1.77	0.50
24:BV:21:ARG:HE	24:BV:87:GLN:HB3	1.75	0.50
1:CA:1338:G:H2'	1:CA:1339:A:C8	2.46	0.50
1:CA:1411:C:H2'	1:CA:1412:C:H6	1.77	0.50
1:CA:676:A:O2'	1:CA:677:U:H5'	2.10	0.50
1:CA:724:G:O2'	1:CA:725:G:H5'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:908:A:H2'	1:CA:909:A:C8	2.46	0.50
3:CD:71:PHE:HA	3:CD:74:TYR:HD2	1.75	0.50
3:CD:81:LEU:HB2	3:CD:88:ASN:ND2	2.25	0.50
4:CE:109:ALA:HB3	4:CE:135:VAL:CG1	2.41	0.50
6:CG:56:SER:HB3	6:CG:59:GLU:HB2	1.92	0.50
9:CJ:66:GLU:HB2	13:CN:98:ALA:CB	2.39	0.50
10:CK:70:ALA:O	10:CK:74:LYS:HB2	2.10	0.50
16:CQ:46:HIS:HB3	16:CQ:73:THR:HG22	1.94	0.50
17:CR:51:GLN:HE21	17:CR:54:LEU:HB2	1.76	0.50
48:D1:7:LYS:HG2	48:D1:26:LYS:N	2.26	0.50
23:DB:1029:A:H2'	23:DB:1030:C:O4'	2.11	0.50
23:DB:1211:C:O2'	23:DB:1212:G:P	2.69	0.50
23:DB:1269:A:H2'	23:DB:1270:C:C6	2.46	0.50
23:DB:1917:U:C2'	23:DB:1918:A:H5'	2.42	0.50
23:DB:1939:U:H5'	23:DB:1939:U:H6	1.75	0.50
23:DB:2076:U:O2	23:DB:2076:U:O4'	2.29	0.50
23:DB:2107:G:H2'	23:DB:2108:A:C8	2.45	0.50
23:DB:2514:U:H2'	23:DB:2515:C:C6	2.47	0.50
23:DB:2899:A:H2'	23:DB:2900:A:H8	1.74	0.50
23:DB:401:A:O2'	23:DB:402:A:H5'	2.11	0.50
23:DB:723:C:H2'	23:DB:724:U:C6	2.47	0.50
23:DB:729:G:H4'	23:DB:763:G:H5'	1.92	0.50
26:DD:127:PHE:C	26:DD:128:ARG:HD3	2.31	0.50
26:DD:37:VAL:HG12	26:DD:44:GLY:O	2.10	0.50
26:DD:70:LYS:HZ2	26:DD:70:LYS:HB3	1.75	0.50
29:DG:120:ILE:HD13	29:DG:140:ILE:HD13	1.93	0.50
30:DH:72:ILE:HG22	30:DH:72:ILE:O	2.11	0.50
52:DI:52:LEU:HD22	52:DI:81:LYS:HD3	1.92	0.50
33:DL:85:VAL:HG21	33:DL:96:LYS:HG3	1.94	0.50
26:DD:116:LYS:HB3	35:DN:1:MET:HA	1.94	0.50
37:DP:28:LYS:O	37:DP:30:TRP:N	2.44	0.50
1:CA:346:G:OP1	37:DP:43:GLU:HG3	2.11	0.50
24:DV:9:ARG:NE	24:DV:20:LEU:HD11	2.27	0.50
43:DW:23:LYS:HG2	43:DW:57:THR:HA	1.92	0.50
1:AA:423:G:H2'	1:AA:424:G:O4'	2.10	0.50
1:AA:685:G:O2'	1:AA:686:U:H5'	2.10	0.50
20:AB:184:ALA:HB3	20:AB:195:VAL:HG21	1.93	0.50
3:AD:96:ARG:O	3:AD:100:VAL:HG23	2.11	0.50
3:AD:197:HIS:O	3:AD:201:GLU:HG3	2.10	0.50
4:AE:71:ILE:HG12	4:AE:72:ASN:N	2.27	0.50
7:AH:110:MET:HB3	7:AH:114:ALA:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AH:40:LYS:HD2	7:AH:48:PHE:CE1	2.47	0.50
8:AI:23:GLY:N	8:AI:60:LEU:HA	2.24	0.50
8:AI:71:ILE:N	8:AI:71:ILE:HD12	2.23	0.50
9:AJ:92:LEU:N	9:AJ:92:LEU:HD22	2.25	0.50
13:AN:30:ILE:CG2	13:AN:41:TRP:HB2	2.37	0.50
15:AP:57:ILE:O	15:AP:61:VAL:HG22	2.11	0.50
16:AQ:35:LYS:O	16:AQ:37:ILE:HG13	2.11	0.50
18:AS:29:PRO:HA	18:AS:47:THR:HB	1.94	0.50
48:B1:10:LEU:HD21	48:B1:34:GLU:HG3	1.93	0.50
22:BA:13:G:O2'	22:BA:15:A:H5''	2.11	0.50
23:BB:1038:G:H2'	23:BB:1039:A:H8	1.76	0.50
23:BB:137:U:H2'	23:BB:138:U:H5'	1.94	0.50
23:BB:2514:U:H2'	23:BB:2515:C:C6	2.46	0.50
23:BB:2630:G:H2'	23:BB:2631:G:H8	1.76	0.50
23:BB:2682:A:O2'	23:BB:2683:C:H5'	2.11	0.50
23:BB:2867:G:N3	23:BB:2867:G:C2'	2.73	0.50
27:BE:165:HIS:O	27:BE:166:LYS:HG3	2.11	0.50
27:BE:46:GLN:O	27:BE:48:THR:N	2.44	0.50
28:BF:116:LEU:H	28:BF:176:PHE:HA	1.76	0.50
29:BG:85:LYS:HA	29:BG:130:ILE:O	2.11	0.50
29:BG:25:ILE:HG23	29:BG:25:ILE:O	2.11	0.50
30:BH:90:LEU:CD2	30:BH:146:VAL:HG11	2.41	0.50
30:BH:68:ARG:HG3	30:BH:134:VAL:CG1	2.38	0.50
52:BI:138:VAL:HG12	52:BI:139:VAL:N	2.26	0.50
52:BI:5:GLN:HG2	52:BI:6:ALA:H	1.74	0.50
23:BB:529:A:OP2	31:BJ:113:PRO:HD3	2.11	0.50
33:BL:83:ALA:HB2	33:BL:120:VAL:HG23	1.94	0.50
23:BB:910:A:C8	34:BM:16:ARG:HB2	2.46	0.50
38:BQ:102:LYS:HG2	38:BQ:103:VAL:N	2.26	0.50
42:BU:23:LYS:HZ3	42:BU:24:VAL:HG22	1.76	0.50
43:BW:24:ARG:HD2	43:BW:57:THR:O	2.12	0.50
44:BX:11:VAL:HG22	44:BX:11:VAL:O	2.10	0.50
44:BX:18:LEU:HD23	44:BX:18:LEU:H	1.76	0.50
1:CA:1520:C:H2'	1:CA:1521:C:H6	1.74	0.50
1:CA:314:C:O2'	1:CA:315:A:H5'	2.11	0.50
1:CA:783:C:O2'	1:CA:784:A:H5'	2.10	0.50
1:CA:853:C:O2'	1:CA:854:U:H5'	2.12	0.50
1:CA:98:A:H2'	1:CA:99:C:H6	1.73	0.50
2:CC:127:VAL:HG23	2:CC:128:MET:N	2.26	0.50
5:CF:39:LEU:HD13	5:CF:39:LEU:O	2.11	0.50
6:CG:45:ALA:HB1	6:CG:120:ALA:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CI:30:ASN:C	8:CI:32:ARG:H	2.14	0.50
1:CA:706:A:H2	10:CK:40:ALA:HB2	1.75	0.50
11:CL:86:VAL:HG11	11:CL:89:LEU:HD23	1.93	0.50
18:CS:44:ILE:HG23	18:CS:63:ASP:HB2	1.94	0.50
18:CS:38:THR:HA	18:CS:69:LYS:HA	1.93	0.50
23:DB:1438:U:H5'	23:DB:1516:G:O2'	2.10	0.50
23:DB:1916:A:H2'	23:DB:1917:U:O4'	2.11	0.50
23:DB:2139:U:O2'	23:DB:2140:G:H5'	2.11	0.50
23:DB:2471:A:O2'	23:DB:2472:G:C8	2.54	0.50
23:DB:2849:U:H4'	23:DB:2868:A:C2	2.46	0.50
23:DB:2831:G:O4'	23:DB:2883:A:C2	2.65	0.50
23:DB:633:A:H2'	23:DB:634:C:C5'	2.41	0.50
23:DB:773:U:H5'	23:DB:774:G:OP2	2.11	0.50
23:DB:850:U:H2'	23:DB:851:C:H6	1.76	0.50
23:DB:1655:A:H4'	26:DD:120:GLY:N	2.26	0.50
27:DE:163:ASN:N	27:DE:168:ASP:HA	2.25	0.50
29:DG:137:LYS:O	29:DG:140:ILE:HB	2.10	0.50
29:DG:23:ILE:HG13	29:DG:23:ILE:O	2.11	0.50
31:DJ:125:TYR:OH	31:DJ:134:ALA:HB2	2.10	0.50
34:DM:127:LYS:NZ	34:DM:129:THR:H	2.08	0.50
38:DQ:60:TRP:CZ3	38:DQ:93:ILE:HG22	2.47	0.50
23:DB:64:A:H5''	41:DT:76:ARG:HG3	1.94	0.50
1:AA:1252:A:H2'	1:AA:1253:G:C5'	2.39	0.50
1:AA:1376:U:H2'	1:AA:1377:A:C8	2.45	0.50
1:AA:1499:A:OP1	53:AA:1601:KSG:N1	2.45	0.50
1:AA:309:A:H2'	1:AA:310:G:H8	1.76	0.50
1:AA:761:G:H2'	1:AA:762:U:C6	2.46	0.50
20:AB:112:ARG:HA	20:AB:115:ASP:OD2	2.11	0.50
2:AC:55:VAL:HG12	2:AC:56:ILE:N	2.26	0.50
3:AD:120:LYS:HB2	3:AD:145:ARG:HH21	1.75	0.50
1:AA:1302:C:O4'	12:AM:16:ILE:HD11	2.12	0.50
12:AM:10:ASP:HA	12:AM:44:ILE:CD1	2.41	0.50
13:AN:30:ILE:C	13:AN:32:ASP:H	2.14	0.50
17:AR:70:THR:HG23	17:AR:71:ASP:N	2.27	0.50
1:AA:957:U:H4'	18:AS:78:THR:HB	1.93	0.50
23:BB:1159:U:O2'	23:BB:1160:G:H5'	2.11	0.50
23:BB:1412:U:H2'	23:BB:1413:A:H8	1.76	0.50
23:BB:1511:G:H2'	23:BB:1512:C:H6	1.77	0.50
23:BB:1676:A:H1'	26:BD:134:HIS:HB3	1.94	0.50
23:BB:2248:C:C2'	23:BB:2249:U:H5'	2.40	0.50
23:BB:2370:G:H2'	23:BB:2371:G:O4'	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2700:A:H2'	23:BB:2701:U:H6	1.76	0.50
23:BB:543:G:H2'	23:BB:544:C:C5'	2.35	0.50
23:BB:550:C:H2'	23:BB:550:C:O2	2.10	0.50
23:BB:771:G:O2'	23:BB:772:C:H5'	2.11	0.50
26:BD:170:VAL:HG11	26:BD:194:PRO:HG2	1.92	0.50
27:BE:44:ARG:N	27:BE:44:ARG:CD	2.73	0.50
29:BG:17:LYS:HD2	29:BG:17:LYS:O	2.09	0.50
52:BI:49:GLU:HG2	52:BI:54:ILE:HD11	1.93	0.50
31:BJ:101:ILE:HG23	31:BJ:102:GLU:N	2.20	0.50
31:BJ:28:LEU:HG	31:BJ:32:LEU:CD1	2.41	0.50
37:BP:109:ILE:O	37:BP:110:LYS:C	2.50	0.50
42:BU:21:ARG:HG2	42:BU:21:ARG:HH11	1.76	0.50
1:CA:114:U:H2'	1:CA:115:G:C8	2.47	0.50
1:CA:1254:A:H61	1:CA:1283:U:H3	1.60	0.50
1:CA:1308:U:H2'	1:CA:1309:G:H8	1.75	0.50
1:CA:1526:G:H2'	1:CA:1527:U:C6	2.46	0.50
1:CA:638:U:H2'	1:CA:639:G:O4'	2.11	0.50
1:CA:846:G:H2'	1:CA:846:G:N3	2.27	0.50
2:CC:69:THR:HG21	2:CC:75:VAL:HG21	1.94	0.50
4:CE:152:VAL:HA	4:CE:155:LYS:HD3	1.93	0.50
10:CK:23:HIS:HB3	10:CK:30:ILE:CG1	2.39	0.50
12:CM:55:LEU:O	12:CM:59:VAL:HG12	2.11	0.50
13:CN:9:GLU:OE2	13:CN:60:ARG:HB3	2.12	0.50
15:CP:36:VAL:HG23	15:CP:56:ARG:HB2	1.93	0.50
19:CT:24:ARG:HA	19:CT:65:LEU:HD21	1.94	0.50
47:D0:26:SER:HB2	47:D0:38:LEU:CG	2.42	0.50
23:DB:145:C:H2'	23:DB:146:A:H8	1.77	0.50
23:DB:1562:U:H2'	23:DB:1563:U:C6	2.46	0.50
23:DB:2665:A:C2'	23:DB:2666:C:H5'	2.41	0.50
23:DB:279:A:H2'	23:DB:280:U:O4'	2.11	0.50
23:DB:2886:A:N7	47:D0:27:LEU:HG	2.26	0.50
23:DB:439:A:O2'	23:DB:440:C:H5'	2.11	0.50
23:DB:490:C:H3'	23:DB:491:G:C5'	2.42	0.50
23:DB:786:C:O2'	23:DB:787:C:H5'	2.10	0.50
25:DC:243:PRO:HA	25:DC:249:VAL:HG23	1.93	0.50
25:DC:92:LEU:HD13	25:DC:102:TYR:CE2	2.47	0.50
23:DB:674:G:O2'	27:DE:69:ARG:HD2	2.12	0.50
29:DG:102:ILE:O	29:DG:102:ILE:HG23	2.12	0.50
30:DH:62:LEU:O	30:DH:66:ASN:HB2	2.12	0.50
31:DJ:131:ASN:C	31:DJ:133:ALA:N	2.64	0.50
33:DL:55:MET:H	33:DL:56:PRO:CD	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DN:33:ILE:HG23	35:DN:112:TYR:HB3	1.93	0.50
35:DN:54:LEU:HD22	35:DN:66:ALA:HB2	1.94	0.50
38:DQ:63:ARG:HG3	38:DQ:96:ASP:OD1	2.12	0.50
40:DS:103:ILE:O	40:DS:104:THR:HB	2.11	0.50
40:DS:72:THR:HG21	40:DS:108:SER:OG	2.11	0.50
41:DT:76:ARG:NH1	41:DT:78:SER:HB2	2.16	0.50
1:AA:1507:A:H2'	1:AA:1508:A:C8	2.47	0.50
1:AA:372:C:H4'	1:AA:373:A:H5'	1.94	0.50
1:AA:602:A:O2'	1:AA:603:U:H5'	2.11	0.50
1:AA:638:U:H2'	1:AA:639:G:O4'	2.11	0.50
1:AA:697:U:H2'	1:AA:698:G:H5'	1.93	0.50
1:AA:751:U:H2'	1:AA:752:G:O4'	2.12	0.50
1:AA:751:U:H4'	14:AO:23:SER:HA	1.92	0.50
1:AA:812:G:O2'	1:AA:813:U:C6	2.58	0.50
20:AB:49:PHE:O	20:AB:53:LEU:HD23	2.11	0.50
2:AC:126:ARG:HH22	2:AC:190:THR:CG2	2.25	0.50
3:AD:24:VAL:C	3:AD:25:ARG:HD2	2.31	0.50
4:AE:81:GLN:HE21	4:AE:147:ASN:C	2.14	0.50
5:AF:99:ALA:O	5:AF:100:SER:HB2	2.11	0.50
10:AK:54:SER:HA	10:AK:56:LYS:HE3	1.92	0.50
15:AP:42:ILE:CG2	15:AP:43:ALA:H	2.18	0.50
47:B0:46:GLY:HA3	47:B0:53:VAL:HG21	1.94	0.50
48:B1:31:GLU:H	48:B1:32:LYS:HZ1	1.59	0.50
49:B2:43:THR:O	49:B2:44:VAL:C	2.49	0.50
23:BB:1019:U:OP1	23:BB:1035:U:O2'	2.29	0.50
23:BB:103:A:H2'	23:BB:104:A:O4'	2.12	0.50
23:BB:1080:A:O2'	23:BB:1081:U:H5'	2.10	0.50
23:BB:1818:U:C3'	25:BC:155:ARG:HB2	2.41	0.50
23:BB:2143:C:H2'	23:BB:2144:G:O4'	2.10	0.50
23:BB:2256:G:O2'	23:BB:2257:U:H5'	2.12	0.50
23:BB:2516:A:O2'	23:BB:2517:C:H5'	2.12	0.50
23:BB:2665:A:C2'	23:BB:2666:C:H5'	2.41	0.50
23:BB:2740:A:H2'	23:BB:2741:A:C8	2.46	0.50
25:BC:155:ARG:HH11	25:BC:157:ALA:HB2	1.76	0.50
26:BD:114:LYS:HB3	35:BN:3:HIS:CE1	2.46	0.50
23:BB:2052:A:O5'	26:BD:147:GLY:HA3	2.12	0.50
26:BD:7:LYS:HE2	26:BD:193:VAL:HG13	1.92	0.50
32:BK:99:ILE:HG22	32:BK:119:ALA:HB2	1.93	0.50
35:BN:17:ARG:O	35:BN:21:PHE:HB2	2.11	0.50
36:BO:18:LEU:HD12	36:BO:18:LEU:C	2.31	0.50
36:BO:54:VAL:HG21	36:BO:78:VAL:HG11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BR:75:VAL:N	39:BR:90:ARG:HE	2.07	0.50
24:BV:76:ASP:HB2	34:BM:136:MET:SD	2.52	0.50
1:CA:370:C:H2'	1:CA:371:A:C8	2.46	0.50
1:CA:761:G:H2'	1:CA:762:U:C6	2.46	0.50
1:CA:963:G:H2'	1:CA:964:A:H8	1.75	0.50
1:CA:973:G:H3'	1:CA:974:A:C5'	2.30	0.50
20:CB:206:ILE:HG22	20:CB:207:ARG:NH2	2.26	0.50
6:CG:16:LYS:HD3	6:CG:16:LYS:O	2.11	0.50
7:CH:17:GLN:HG3	7:CH:71:VAL:CG2	2.41	0.50
13:CN:13:VAL:HG13	13:CN:59:GLN:HE22	1.76	0.50
18:CS:52:ASN:ND2	18:CS:76:THR:HA	2.26	0.50
22:DA:31:C:O2'	22:DA:53:A:N1	2.44	0.50
23:DB:1396:U:O4'	23:DB:1396:U:O2	2.28	0.50
23:DB:1681:G:H2'	23:DB:1757:A:N1	2.27	0.50
23:DB:2056:G:H21	47:D0:1:ALA:HB3	1.76	0.50
23:DB:2412:A:H2'	23:DB:2413:G:O4'	2.11	0.50
23:DB:2636:C:H2'	23:DB:2637:U:C6	2.46	0.50
23:DB:2743:U:C3'	23:DB:2744:G:H5''	2.42	0.50
23:DB:363:G:H2'	23:DB:364:C:H6	1.74	0.50
23:DB:910:A:C8	34:DM:16:ARG:HB3	2.47	0.50
25:DC:103:ILE:HG22	25:DC:104:LEU:H	1.77	0.50
25:DC:207:ALA:HA	25:DC:211:ARG:HB3	1.94	0.50
25:DC:3:VAL:HG12	25:DC:4:LYS:N	2.27	0.50
26:DD:117:GLY:O	26:DD:164:GLN:HA	2.12	0.50
26:DD:60:VAL:O	26:DD:63:PRO:HD2	2.12	0.50
23:DB:2636:C:P	26:DD:80:TRP:HE1	2.34	0.50
27:DE:192:ALA:HB1	27:DE:199:MET:CG	2.42	0.50
52:DI:99:LYS:HD3	52:DI:99:LYS:N	2.26	0.50
31:DJ:37:ARG:NH1	31:DJ:110:PRO:HG3	2.27	0.50
31:DJ:5:THR:HG21	31:DJ:7:LYS:NZ	2.26	0.50
32:DK:110:GLU:HA	32:DK:113:MET:HE3	1.92	0.50
33:DL:118:THR:CG2	33:DL:137:ALA:HB3	2.41	0.50
33:DL:58:TYR:HA	33:DL:62:PRO:CD	2.42	0.50
35:DN:115:LEU:N	35:DN:115:LEU:HD23	2.26	0.50
36:DO:80:GLU:O	36:DO:83:LEU:HD13	2.10	0.50
37:DP:100:ARG:H	37:DP:100:ARG:CD	2.25	0.50
38:DQ:5:ARG:HD2	38:DQ:5:ARG:O	2.11	0.50
40:DS:64:ALA:H	40:DS:110:ARG:NH2	2.10	0.50
40:DS:42:LYS:HD3	40:DS:42:LYS:O	2.11	0.50
43:DW:48:ALA:O	43:DW:72:GLY:HA3	2.12	0.50
1:AA:1004:A:H62	1:AA:1025:U:H5'	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1210:C:O2'	1:AA:1211:U:H5'	2.11	0.50
1:AA:1504:G:H4'	1:AA:1505:G:C4	2.47	0.50
1:AA:437:U:C2'	1:AA:438:U:O4'	2.51	0.50
1:AA:60:A:H1'	1:AA:61:G:O4'	2.12	0.50
1:AA:769:G:O2'	1:AA:770:C:H5'	2.11	0.50
1:AA:864:A:H2'	1:AA:865:A:C8	2.46	0.50
2:AC:183:TYR:HE1	2:AC:198:LYS:HB3	1.75	0.50
2:AC:71:ARG:O	2:AC:75:VAL:HG23	2.11	0.50
5:AF:45:ARG:HG2	5:AF:46:GLN:N	2.26	0.50
6:AG:46:LEU:O	6:AG:46:LEU:HD13	2.11	0.50
8:AI:11:ARG:HG3	8:AI:11:ARG:O	2.10	0.50
50:B3:7:ARG:HH21	50:B3:11:LYS:HZ2	1.57	0.50
23:BB:1373:A:OP1	23:BB:2213:U:O4	2.29	0.50
23:BB:1531:C:H2'	23:BB:1532:A:H8	1.75	0.50
23:BB:1796:U:H2'	23:BB:1797:G:C8	2.46	0.50
23:BB:2025:C:H2'	23:BB:2026:U:H6	1.77	0.50
23:BB:2262:U:H4'	23:BB:2328:A:H2	1.76	0.50
23:BB:2334:U:OP2	36:BO:7:ARG:HA	2.11	0.50
23:BB:2439:A:C8	23:BB:2586:U:H4'	2.47	0.50
23:BB:408:G:O2'	23:BB:409:G:H5'	2.12	0.50
23:BB:557:C:H2'	23:BB:558:U:H6	1.76	0.50
23:BB:833:A:H2'	23:BB:834:G:H8	1.74	0.50
25:BC:61:TYR:CE2	25:BC:63:ILE:HD11	2.47	0.50
27:BE:131:THR:O	27:BE:134:LEU:HB2	2.12	0.50
27:BE:139:LYS:HE2	27:BE:139:LYS:HA	1.93	0.50
28:BF:168:LEU:HD12	28:BF:169:LEU:N	2.15	0.50
29:BG:18:ILE:CD1	29:BG:18:ILE:H	2.25	0.50
30:BH:58:LEU:HA	30:BH:62:LEU:HD12	1.94	0.50
32:BK:23:LYS:HG3	32:BK:24:VAL:H	1.75	0.50
33:BL:109:LYS:HA	33:BL:126:ARG:CB	2.42	0.50
33:BL:123:ARG:HH22	33:BL:141:LYS:HZ1	1.57	0.50
36:BO:31:THR:HG21	36:BO:36:TYR:CE1	2.46	0.50
38:BQ:40:LYS:HA	38:BQ:43:GLN:HE21	1.77	0.50
42:BU:82:VAL:HB	42:BU:94:PHE:CD2	2.47	0.50
1:CA:1008:U:H5''	13:CN:23:ARG:NH2	2.20	0.50
1:CA:1118:U:H5'	1:CA:1118:U:H6	1.76	0.50
1:CA:1200:C:H4'	1:CA:1201:A:H5'	1.94	0.50
1:CA:1427:C:O2'	1:CA:1428:A:H5'	2.11	0.50
1:CA:143:A:H2	1:CA:220:G:H22	1.59	0.50
1:CA:1492:A:N6	1:CA:1494:G:C8	2.80	0.50
20:CB:115:ASP:O	20:CB:119:GLN:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:169:TRP:CE3	3:CD:185:PRO:HB3	2.46	0.50
4:CE:15:ILE:HG22	4:CE:16:ALA:H	1.77	0.50
4:CE:52:ALA:HB2	4:CE:61:LYS:HE2	1.93	0.50
5:CF:29:ILE:HG21	5:CF:64:VAL:HG11	1.93	0.50
9:CJ:66:GLU:HG2	9:CJ:67:ILE:H	1.77	0.50
13:CN:40:ARG:O	13:CN:45:LEU:HD11	2.11	0.50
13:CN:42:ASN:HA	13:CN:45:LEU:HD12	1.94	0.50
15:CP:25:ARG:HD3	15:CP:25:ARG:H	1.76	0.50
1:CA:1221:G:OP1	18:CS:35:ARG:HD2	2.11	0.50
51:D4:22:VAL:HG13	51:D4:37:GLN:HB3	1.93	0.50
22:DA:66:A:N6	22:DA:107:G:C5	2.79	0.50
23:DB:1313:U:H4'	23:DB:1332:G:H4'	1.94	0.50
23:DB:2020:A:C2	23:DB:2022:U:O4'	2.64	0.50
23:DB:2405:G:H1'	23:DB:2412:A:H61	1.75	0.50
23:DB:2742:G:O2'	23:DB:2743:U:H5'	2.12	0.50
23:DB:2771:C:H2'	23:DB:2772:C:H6	1.76	0.50
23:DB:46:G:H2'	23:DB:47:C:H6	1.77	0.50
23:DB:935:C:H2'	23:DB:936:A:C8	2.47	0.50
23:DB:969:G:OP1	45:DY:17:PRO:HG3	2.11	0.50
25:DC:216:ARG:O	25:DC:218:THR:N	2.44	0.50
25:DC:268:ARG:HG3	25:DC:268:ARG:HH11	1.77	0.50
26:DD:146:ILE:CD1	26:DD:146:ILE:H	2.24	0.50
23:DB:1205:A:C8	27:DE:165:HIS:HA	2.47	0.50
27:DE:142:ALA:N	27:DE:185:LYS:NZ	2.52	0.50
28:DF:98:PHE:HA	28:DF:101:ARG:HG2	1.93	0.50
30:DH:122:LEU:HD13	30:DH:146:VAL:CG2	2.36	0.50
31:DJ:133:ALA:C	31:DJ:135:GLN:N	2.65	0.50
32:DK:11:ALA:O	32:DK:99:ILE:HG23	2.10	0.50
35:DN:10:LEU:O	35:DN:11:ASN:ND2	2.45	0.50
35:DN:73:ASN:HA	35:DN:76:VAL:HG12	1.93	0.50
37:DP:25:VAL:O	37:DP:27:VAL:HG12	2.11	0.50
37:DP:55:HIS:C	37:DP:57:ALA:N	2.64	0.50
39:DR:47:VAL:HG22	39:DR:48:LYS:N	2.18	0.50
41:DT:53:VAL:HB	41:DT:93:LEU:HD21	1.93	0.50
44:DX:16:THR:OG1	44:DX:19:LEU:HB3	2.12	0.50
1:AA:1261:A:H2'	1:AA:1262:C:O4'	2.11	0.50
1:AA:1324:A:H2'	1:AA:1325:C:C6	2.47	0.50
1:AA:146:G:O2'	1:AA:147:G:H5'	2.11	0.50
1:AA:251:G:H4'	1:AA:252:U:H5'	1.93	0.50
1:AA:323:U:H2'	1:AA:324:G:O4'	2.11	0.50
1:AA:980:C:H2'	1:AA:981:U:H5'	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AB:80:LYS:O	20:AB:84:LEU:N	2.45	0.50
2:AC:155:ARG:NH1	2:AC:192:TYR:HB2	2.27	0.50
4:AE:101:GLY:H	4:AE:121:ASN:ND2	2.07	0.50
7:AH:76:ARG:HE	7:AH:79:ARG:HA	1.77	0.50
8:AI:50:PRO:HG2	8:AI:51:LEU:HD12	1.92	0.50
21:AU:26:GLY:C	21:AU:28:LEU:H	2.15	0.50
47:B0:11:LYS:O	47:B0:12:ARG:HB3	2.12	0.50
48:B1:36:LYS:N	48:B1:36:LYS:HD2	2.27	0.50
22:BA:32:U:H4'	22:BA:52:A:H62	1.76	0.50
23:BB:1236:G:O2'	23:BB:1237:A:O5'	2.29	0.50
23:BB:1313:U:H4'	23:BB:1332:G:H4'	1.94	0.50
23:BB:1396:U:O2	23:BB:1396:U:O4'	2.28	0.50
23:BB:1909:C:H2'	23:BB:1910:G:H8	1.77	0.50
23:BB:2020:A:C2	23:BB:2022:U:O4'	2.64	0.50
23:BB:235:U:H2'	23:BB:236:C:H6	1.76	0.50
23:BB:2669:G:H2'	23:BB:2670:A:H8	1.76	0.50
23:BB:2821:A:H2'	23:BB:2822:G:C8	2.46	0.50
23:BB:38:A:C8	27:BE:46:GLN:CD	2.85	0.50
23:BB:536:G:H4'	38:BQ:56:PHE:HZ	1.74	0.50
23:BB:855:G:N2	43:BW:23:LYS:HB3	2.26	0.50
23:BB:993:G:O2'	23:BB:994:C:H5'	2.11	0.50
25:BC:136:VAL:O	25:BC:163:ILE:HB	2.12	0.50
25:BC:67:LYS:HB3	25:BC:149:LYS:O	2.12	0.50
25:BC:77:VAL:HA	25:BC:93:VAL:HG22	1.94	0.50
26:BD:149:ASN:C	26:BD:150:GLN:HG2	2.32	0.50
23:BB:2822:G:OP1	26:BD:166:GLY:HA2	2.11	0.50
27:BE:104:ALA:O	27:BE:105:LEU:HD23	2.12	0.50
27:BE:170:ARG:NH2	27:BE:175:ILE:HA	2.26	0.50
27:BE:57:LYS:O	27:BE:59:PRO:HD3	2.11	0.50
28:BF:106:ALA:CB	28:BF:136:ILE:HG23	2.42	0.50
30:BH:85:GLY:O	30:BH:89:LYS:N	2.45	0.50
32:BK:10:VAL:HG13	32:BK:19:VAL:HG21	1.94	0.50
23:BB:2675:A:OP1	32:BK:31:ARG:HG2	2.10	0.50
33:BL:71:ALA:O	33:BL:72:ALA:C	2.50	0.50
33:BL:92:LEU:HG	33:BL:96:LYS:NZ	2.27	0.50
34:BM:26:VAL:HG11	34:BM:134:THR:OG1	2.12	0.50
23:BB:2485:G:H5'	34:BM:45:GLN:OE1	2.12	0.50
36:BO:59:ALA:HB3	36:BO:63:LYS:HB2	1.93	0.50
37:BP:61:ARG:O	37:BP:63:ILE:HD12	2.11	0.50
38:BQ:111:LYS:O	38:BQ:115:ALA:HB2	2.12	0.50
38:BQ:13:HIS:O	38:BQ:17:LEU:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BQ:7:VAL:HG23	38:BQ:8:ILE:N	2.18	0.50
38:BQ:88:GLU:HA	39:BR:53:PHE:CD1	2.47	0.50
42:BU:24:VAL:C	42:BU:25:LYS:HD3	2.32	0.50
23:BB:988:A:OP2	45:BY:11:SER:HA	2.12	0.50
1:CA:1026:G:H2'	1:CA:1027:C:O4'	2.12	0.50
1:CA:1171:A:H2'	1:CA:1172:C:H6	1.77	0.50
1:CA:1219:A:H2'	1:CA:1220:G:H8	1.75	0.50
1:CA:1496:C:H2'	1:CA:1497:G:O4'	2.11	0.50
1:CA:1512:U:H2'	1:CA:1513:A:H8	1.75	0.50
1:CA:806:C:H2'	1:CA:807:A:C8	2.46	0.50
20:CB:85:SER:O	20:CB:86:CYS:HB2	2.12	0.50
3:CD:84:ASN:HD21	3:CD:87:GLU:H	1.58	0.50
5:CF:68:GLN:O	5:CF:71:ILE:HG22	2.11	0.50
6:CG:29:LEU:HD11	6:CG:119:LEU:HD22	1.94	0.50
7:CH:77:VAL:CG2	7:CH:126:CYS:HA	2.42	0.50
12:CM:102:LYS:HG3	12:CM:103:THR:N	2.26	0.50
16:CQ:43:LEU:HD12	16:CQ:43:LEU:N	2.27	0.50
23:DB:2399:G:H1'	48:D1:20:TYR:OH	2.12	0.50
23:DB:2420:C:H3'	50:D3:31:ILE:CG2	2.42	0.50
23:DB:1139:G:O2'	23:DB:1140:C:H5'	2.12	0.50
23:DB:125:A:O2'	23:DB:126:A:OP1	2.24	0.50
23:DB:1948:G:O2'	23:DB:1949:G:H5'	2.12	0.50
23:DB:2464:G:H2'	23:DB:2465:C:H6	1.75	0.50
23:DB:2500:U:H5'	23:DB:2501:C:OP2	2.12	0.50
23:DB:242:G:H22	23:DB:254:G:H2'	1.76	0.50
23:DB:299:A:H2'	23:DB:300:A:C8	2.46	0.50
23:DB:441:U:H2'	23:DB:442:G:H8	1.76	0.50
23:DB:725:G:H2'	23:DB:726:G:C1'	2.41	0.50
23:DB:943:A:OP1	33:DL:41:ARG:HG2	2.12	0.50
25:DC:21:PRO:C	25:DC:202:ARG:HD2	2.32	0.50
27:DE:199:MET:O	27:DE:200:LEU:HD23	2.12	0.50
27:DE:29:HIS:HD2	33:DL:19:LEU:HD11	1.76	0.50
28:DF:174:PHE:N	28:DF:175:PRO:HD2	2.23	0.50
52:DI:21:PRO:CB	52:DI:22:PRO:HD3	2.39	0.50
52:DI:85:ILE:HD12	52:DI:87:SER:O	2.12	0.50
34:DM:16:ARG:HG2	34:DM:16:ARG:HH11	1.77	0.50
41:DT:11:LEU:HB3	41:DT:34:VAL:HG23	1.92	0.50
24:DV:57:TYR:HE2	24:DV:77:VAL:HG21	1.76	0.50
24:DV:65:VAL:C	24:DV:67:GLY:H	2.15	0.50
43:DW:42:THR:H	43:DW:65:LYS:CG	2.22	0.50
1:AA:1226:C:H4'	1:AA:1227:A:OP1	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:143:A:H2	1:AA:220:G:H22	1.59	0.50
1:AA:219:U:H2'	1:AA:220:G:H8	1.77	0.50
1:AA:370:C:H2'	1:AA:371:A:H8	1.77	0.50
1:AA:969:A:N3	1:AA:970:C:O2	2.45	0.50
20:AB:206:ILE:HD13	20:AB:206:ILE:C	2.32	0.50
2:AC:180:ASP:C	2:AC:181:ILE:HD12	2.31	0.50
3:AD:11:SER:OG	3:AD:17:ASP:HA	2.11	0.50
3:AD:77:GLU:O	3:AD:81:LEU:HG	2.12	0.50
5:AF:51:ILE:HD11	5:AF:86:ARG:HG3	1.94	0.50
8:AI:49:GLN:N	8:AI:50:PRO:HD2	2.27	0.50
8:AI:56:MET:HE2	8:AI:57:VAL:H	1.77	0.50
1:AA:130:A:O4'	16:AQ:64:ARG:HD2	2.12	0.50
1:AA:734:G:N2	17:AR:63:TYR:CE2	2.80	0.50
22:BA:52:A:H3'	22:BA:53:A:C8	2.47	0.50
23:BB:1386:C:H2'	23:BB:1387:A:H8	1.74	0.50
23:BB:2809:A:H2'	23:BB:2810:A:H8	1.77	0.50
23:BB:827:U:H5'	23:BB:828:U:O5'	2.12	0.50
23:BB:922:C:H2'	23:BB:923:G:C8	2.47	0.50
23:BB:968:C:H2'	23:BB:969:G:C8	2.45	0.50
25:BC:130:PRO:HA	25:BC:187:CYS:O	2.12	0.50
25:BC:22:GLU:HA	25:BC:202:ARG:NE	2.26	0.50
26:BD:45:TYR:O	26:BD:46:ARG:HG2	2.12	0.50
26:BD:48:ILE:HA	26:BD:79:LEU:O	2.10	0.50
27:BE:120:VAL:HG12	27:BE:121:VAL:N	2.27	0.50
27:BE:88:ARG:HH11	27:BE:88:ARG:CB	2.22	0.50
29:BG:102:ILE:HG22	29:BG:114:HIS:O	2.12	0.50
30:BH:69:ALA:HB1	30:BH:141:LYS:NZ	2.27	0.50
31:BJ:7:LYS:CD	31:BJ:48:VAL:HB	2.39	0.50
31:BJ:61:LYS:CE	31:BJ:61:LYS:HA	2.38	0.50
32:BK:16:ALA:N	32:BK:46:ALA:HA	2.27	0.50
33:BL:116:VAL:HG21	33:BL:134:ALA:HB1	1.94	0.50
35:BN:1:MET:SD	35:BN:2:ARG:N	2.85	0.50
37:BP:61:ARG:NH2	37:BP:63:ILE:HD11	2.27	0.50
23:BB:535:G:N2	38:BQ:52:ARG:HH22	1.95	0.50
38:BQ:56:PHE:O	38:BQ:59:LEU:HB3	2.12	0.50
38:BQ:94:LEU:HD11	39:BR:10:LYS:HD2	1.94	0.50
39:BR:49:ILE:HG22	39:BR:51:VAL:CG2	2.42	0.50
40:BS:34:ASP:HA	40:BS:38:TYR:CD2	2.47	0.50
41:BT:28:ASN:HD22	41:BT:28:ASN:N	2.09	0.50
43:BW:4:LYS:C	43:BW:4:LYS:HD3	2.32	0.50
45:BY:6:ILE:HG22	45:BY:56:VAL:CG2	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:BY:9:THR:HA	45:BY:55:LYS:HG3	1.92	0.50
1:CA:1307:U:H2'	1:CA:1308:U:H6	1.76	0.50
3:CD:138:PRO:O	3:CD:139:ASN:HB3	2.12	0.50
4:CE:111:ARG:HG2	4:CE:111:ARG:NH1	2.27	0.50
9:CJ:85:ASP:HB2	9:CJ:89:ARG:CZ	2.42	0.50
10:CK:108:ASN:OD1	10:CK:110:THR:HG23	2.12	0.50
11:CL:113:ARG:CZ	11:CL:120:ARG:HA	2.42	0.50
15:CP:74:LEU:O	15:CP:78:VAL:HG12	2.11	0.50
5:CF:100:SER:H	17:CR:23:LYS:HZ1	1.60	0.50
18:CS:20:LYS:HD2	18:CS:20:LYS:O	2.12	0.50
13:CN:40:ARG:HH12	18:CS:6:LYS:H	1.59	0.50
22:DA:30:C:C2'	22:DA:31:C:H5'	2.39	0.50
23:DB:1184:U:O2'	23:DB:1185:G:H5'	2.12	0.50
23:DB:1330:C:H2'	23:DB:1331:G:H8	1.76	0.50
23:DB:699:A:H4'	23:DB:1634:A:N7	2.27	0.50
23:DB:1771:C:H2'	23:DB:1772:A:C8	2.46	0.50
23:DB:2052:A:OP1	26:DD:145:SER:HB3	2.12	0.50
23:DB:2520:C:C6	23:DB:2567:G:H1'	2.46	0.50
23:DB:2630:G:H2'	23:DB:2631:G:H8	1.75	0.50
23:DB:678:C:H2'	23:DB:679:C:H6	1.77	0.50
25:DC:192:GLY:O	25:DC:194:VAL:HG22	2.11	0.50
26:DD:160:LYS:HA	26:DD:160:LYS:HE3	1.93	0.50
28:DF:134:GLN:O	28:DF:135:ILE:C	2.48	0.50
28:DF:137:PHE:O	28:DF:139:GLU:HG2	2.12	0.50
52:DI:78:LEU:HD23	52:DI:81:LYS:HE2	1.93	0.50
33:DL:51:GLU:CG	33:DL:52:GLY:H	2.09	0.50
34:DM:34:LYS:O	34:DM:125:PRO:HD2	2.12	0.50
26:DD:116:LYS:O	35:DN:2:ARG:HB3	2.12	0.50
36:DO:95:SER:O	36:DO:97:PHE:N	2.45	0.50
43:DW:43:LYS:H	43:DW:75:ASN:CG	2.15	0.50
44:DX:25:GLN:CD	44:DX:29:ARG:HH21	2.14	0.50
1:AA:1446:A:C3'	1:AA:1447:A:H5''	2.42	0.50
1:AA:193:C:H2'	1:AA:194:C:C6	2.47	0.50
1:AA:939:G:H4'	6:AG:101:ARG:NH2	2.26	0.50
6:AG:136:LYS:O	6:AG:140:VAL:HG23	2.12	0.50
6:AG:72:VAL:HB	6:AG:144:ALA:HB3	1.94	0.50
8:AI:56:MET:CE	8:AI:57:VAL:H	2.25	0.50
14:AO:70:LYS:NZ	14:AO:74:VAL:HA	2.27	0.50
15:AP:67:ILE:HG13	15:AP:71:VAL:HG13	1.93	0.50
16:AQ:80:LYS:C	16:AQ:80:LYS:HE3	2.32	0.50
19:AT:15:LYS:HA	19:AT:18:LYS:HE3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AU:34:ARG:C	21:AU:34:ARG:HD2	2.31	0.50
21:AU:5:VAL:HG22	21:AU:19:LYS:NZ	2.27	0.50
23:BB:1107:G:H2'	23:BB:1108:U:O4'	2.11	0.50
23:BB:1399:C:H2'	23:BB:1400:U:C6	2.47	0.50
23:BB:189:G:P	46:BZ:12:ILE:HG12	2.52	0.50
23:BB:2087:G:H2'	23:BB:2088:A:H8	1.77	0.50
23:BB:2877:G:H2'	23:BB:2878:U:C6	2.47	0.50
23:BB:2811:G:H5''	26:BD:62:LYS:HD2	1.93	0.50
27:BE:4:VAL:HB	27:BE:117:ARG:HH21	1.76	0.50
27:BE:50:ALA:O	27:BE:74:LYS:HD2	2.12	0.50
23:BB:2305:U:H3	28:BF:150:GLY:HA3	1.76	0.50
30:BH:6:LEU:HD12	30:BH:15:LEU:CA	2.42	0.50
30:BH:57:LYS:CA	30:BH:57:LYS:HE3	2.42	0.50
30:BH:56:ALA:HB3	30:BH:57:LYS:NZ	2.27	0.50
31:BJ:18:VAL:HG22	31:BJ:56:VAL:HA	1.93	0.50
31:BJ:36:LEU:O	31:BJ:118:MET:HG2	2.12	0.50
34:BM:112:LEU:N	34:BM:112:LEU:HD12	2.27	0.50
34:BM:97:GLN:N	34:BM:98:PRO:CD	2.75	0.50
36:BO:17:LYS:O	36:BO:21:LEU:HA	2.12	0.50
40:BS:17:VAL:HA	40:BS:43:ALA:CB	2.42	0.50
40:BS:57:ASN:O	40:BS:61:ASN:HB3	2.12	0.50
41:BT:13:ALA:HB3	41:BT:33:LYS:HB2	1.92	0.50
24:BV:80:HIS:CD2	24:BV:82:TYR:H	2.25	0.50
44:BX:5:GLU:N	44:BX:7:ARG:HE	2.08	0.50
46:BZ:49:ARG:CD	46:BZ:50:ASP:N	2.75	0.50
1:CA:1129:C:H1'	1:CA:1146:A:N6	2.25	0.50
1:CA:262:A:OP2	19:CT:70:LYS:HE3	2.12	0.50
1:CA:372:C:H4'	1:CA:373:A:H5'	1.94	0.50
1:CA:736:C:H2'	1:CA:737:C:H6	1.77	0.50
1:CA:945:G:H21	1:CA:1334:G:H4'	1.77	0.50
20:CB:134:LEU:O	20:CB:134:LEU:HD23	2.11	0.50
15:CP:67:ILE:CG1	15:CP:71:VAL:HG13	2.37	0.50
16:CQ:13:SER:HB3	16:CQ:21:VAL:CB	2.40	0.50
18:CS:11:ASP:HB3	18:CS:13:HIS:NE2	2.27	0.50
18:CS:16:LYS:O	18:CS:20:LYS:HB2	2.11	0.50
21:CU:3:ILE:HD12	21:CU:3:ILE:O	2.11	0.50
51:D4:2:LYS:HD3	51:D4:2:LYS:N	2.27	0.50
23:DB:1196:C:H2'	23:DB:1197:G:C8	2.46	0.50
23:DB:19:A:OP1	38:DQ:22:GLY:HA2	2.12	0.50
23:DB:217:A:H3'	23:DB:218:A:H8	1.75	0.50
23:DB:2191:A:H2'	23:DB:2192:U:C6	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2682:A:O2'	23:DB:2683:C:H5'	2.11	0.50
23:DB:2877:G:H2'	23:DB:2878:U:C6	2.47	0.50
23:DB:416:U:H2'	23:DB:417:C:C6	2.46	0.50
23:DB:553:G:O2'	23:DB:554:U:H5'	2.11	0.50
27:DE:132:LYS:HG2	27:DE:134:LEU:CD1	2.41	0.50
27:DE:164:LEU:HD13	27:DE:165:HIS:N	2.27	0.50
27:DE:28:VAL:O	27:DE:32:VAL:HG23	2.12	0.50
29:DG:51:PHE:CE1	29:DG:53:PRO:HG3	2.45	0.50
30:DH:116:ARG:O	30:DH:116:ARG:HG2	2.12	0.50
30:DH:114:GLU:CB	30:DH:133:GLN:HG3	2.37	0.50
31:DJ:78:THR:H	31:DJ:84:ILE:CD1	2.25	0.50
31:DJ:76:HIS:HB2	31:DJ:86:GLN:CG	2.41	0.50
33:DL:35:HIS:CG	33:DL:35:HIS:O	2.65	0.50
33:DL:44:GLY:HA2	33:DL:47:ARG:NE	2.22	0.50
36:DO:40:ILE:N	36:DO:40:ILE:HD13	2.16	0.50
39:DR:72:VAL:O	39:DR:72:VAL:HG13	2.11	0.50
44:DX:18:LEU:HD12	44:DX:47:ARG:HH22	1.77	0.50
1:AA:1169:A:H2'	1:AA:1170:A:C8	2.48	0.49
1:AA:1332:A:H2'	1:AA:1333:A:C8	2.47	0.49
1:AA:1458:G:H2'	1:AA:1459:G:H8	1.77	0.49
1:AA:33:A:H2'	1:AA:34:C:H6	1.77	0.49
1:AA:968:A:O2'	1:AA:969:A:OP1	2.27	0.49
20:AB:46:VAL:O	20:AB:49:PHE:HB2	2.12	0.49
11:AL:20:VAL:HG12	11:AL:20:VAL:O	2.12	0.49
16:AQ:10:ARG:NH2	16:AQ:58:VAL:HG13	2.27	0.49
16:AQ:66:LEU:HD11	16:AQ:73:THR:CG2	2.41	0.49
17:AR:44:THR:HG22	17:AR:46:THR:HB	1.95	0.49
21:AU:26:GLY:O	21:AU:30:GLU:HB2	2.12	0.49
47:B0:46:GLY:O	47:B0:52:LYS:HB3	2.12	0.49
23:BB:1054:A:H2'	23:BB:1055:G:C8	2.45	0.49
23:BB:2526:G:H2'	23:BB:2527:C:C6	2.47	0.49
23:BB:272:A:H2'	23:BB:273:G:C8	2.47	0.49
23:BB:395:U:H2'	23:BB:396:G:C8	2.47	0.49
23:BB:708:G:H2'	23:BB:709:U:C6	2.47	0.49
23:BB:932:U:H2'	23:BB:934:U:C4	2.46	0.49
25:BC:172:THR:HG22	25:BC:173:LEU:N	2.24	0.49
25:BC:222:THR:HG23	25:BC:231:HIS:CD2	2.47	0.49
25:BC:182:LYS:HE2	25:BC:264:LYS:NZ	2.27	0.49
26:BD:8:LYS:HB3	37:BP:5:LYS:HZ2	1.74	0.49
28:BF:71:LYS:HD3	28:BF:71:LYS:C	2.32	0.49
29:BG:126:THR:HG23	29:BG:128:THR:HG23	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:BH:5:LEU:HD21	30:BH:17:ASP:OD1	2.12	0.49
30:BH:48:GLU:C	30:BH:50:ARG:H	2.15	0.49
32:BK:23:LYS:O	32:BK:39:ILE:HG23	2.12	0.49
26:BD:18:ASP:O	32:BK:73:ASP:HA	2.11	0.49
33:BL:18:ARG:N	33:BL:18:ARG:HD3	2.27	0.49
34:BM:38:ARG:HA	34:BM:96:ILE:O	2.11	0.49
37:BP:38:ARG:NH1	37:BP:39:LEU:N	2.59	0.49
38:BQ:97:ILE:HD13	38:BQ:100:PHE:CZ	2.46	0.49
38:BQ:90:ASP:HA	39:BR:10:LYS:HZ3	1.76	0.49
42:BU:15:GLY:HA2	42:BU:18:LYS:HD3	1.92	0.49
1:CA:1070:U:H2'	1:CA:1071:C:C6	2.47	0.49
1:CA:1241:G:H2'	1:CA:1242:G:C8	2.42	0.49
1:CA:1298:U:H2'	6:CG:113:LYS:NZ	2.27	0.49
1:CA:1492:A:H2'	1:CA:1492:A:N3	2.26	0.49
1:CA:22:G:H2'	1:CA:23:C:C6	2.47	0.49
1:CA:309:A:H2'	1:CA:310:G:H8	1.76	0.49
1:CA:751:U:H2'	1:CA:752:G:O4'	2.12	0.49
20:CB:45:THR:HG23	20:CB:200:PRO:HG2	1.94	0.49
20:CB:212:TYR:O	20:CB:216:VAL:HG22	2.11	0.49
1:CA:829:G:H4'	20:CB:24:PRO:HG3	1.94	0.49
2:CC:149:LYS:HE3	2:CC:200:TRP:CE3	2.47	0.49
2:CC:61:LYS:NZ	2:CC:96:VAL:HG12	2.26	0.49
1:CA:429:U:P	3:CD:12:ARG:HH21	2.35	0.49
6:CG:55:LYS:HB3	6:CG:59:GLU:OE2	2.12	0.49
10:CK:85:VAL:HG11	10:CK:96:ILE:HD11	1.93	0.49
10:CK:22:ILE:HG21	10:CK:95:THR:HG21	1.93	0.49
14:CO:72:LYS:O	14:CO:73:ASP:HB2	2.11	0.49
18:CS:10:ILE:HD13	18:CS:40:PHE:CE2	2.47	0.49
48:D1:9:LYS:HG2	48:D1:24:LYS:HG2	1.93	0.49
49:D2:12:ARG:CG	49:D2:46:LYS:HA	2.42	0.49
48:D1:22:THR:HG21	50:D3:34:LYS:HZ1	1.77	0.49
23:DB:1399:C:H2'	23:DB:1400:U:C6	2.47	0.49
23:DB:1439:A:N7	23:DB:1440:U:C2	2.80	0.49
23:DB:1501:G:O2'	23:DB:1502:A:H5'	2.11	0.49
23:DB:1533:C:C2'	23:DB:1534:U:H5'	2.42	0.49
23:DB:1796:U:H2'	23:DB:1797:G:C8	2.46	0.49
23:DB:2106:U:H2'	23:DB:2107:G:OP1	2.12	0.49
23:DB:2514:U:H2'	23:DB:2515:C:H6	1.77	0.49
23:DB:2439:A:C8	23:DB:2586:U:H4'	2.47	0.49
23:DB:351:C:H2'	23:DB:352:A:C8	2.46	0.49
23:DB:1813:G:C1'	25:DC:45:ASN:HB3	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:193:VAL:HG23	26:DD:193:VAL:O	2.11	0.49
27:DE:130:LYS:HZ3	27:DE:130:LYS:HB2	1.77	0.49
27:DE:193:VAL:HG12	27:DE:199:MET:H	1.76	0.49
28:DF:107:VAL:N	28:DF:108:PRO:HD2	2.26	0.49
28:DF:115:GLY:HA3	28:DF:177:ARG:HB2	1.94	0.49
28:DF:41:GLU:O	28:DF:42:ALA:C	2.50	0.49
29:DG:53:PRO:HG2	29:DG:61:TRP:HZ3	1.72	0.49
23:DB:1080:A:H4'	52:DI:126:ARG:HD3	1.93	0.49
32:DK:53:LYS:O	32:DK:56:ASP:HB2	2.12	0.49
35:DN:58:ASP:O	35:DN:62:ASN:HB2	2.10	0.49
42:DU:10:VAL:O	42:DU:21:ARG:HA	2.12	0.49
44:DX:17:GLU:H	44:DX:17:GLU:CD	2.14	0.49
1:AA:1363:A:N3	1:AA:1363:A:H2'	2.27	0.49
1:AA:189:A:H2'	1:AA:190:A:C8	2.47	0.49
1:AA:399:G:H2'	1:AA:400:C:H6	1.77	0.49
20:AB:111:LYS:O	20:AB:114:LYS:HB2	2.11	0.49
20:AB:202:ASN:HD22	20:AB:204:ASP:N	1.99	0.49
20:AB:41:ASN:ND2	20:AB:44:LYS:HE2	2.26	0.49
2:AC:151:GLU:HB2	2:AC:200:TRP:HZ3	1.75	0.49
5:AF:38:ARG:HH21	5:AF:96:VAL:HG12	1.77	0.49
7:AH:49:LYS:HG3	7:AH:50:VAL:N	2.25	0.49
7:AH:98:LEU:HD12	7:AH:98:LEU:H	1.77	0.49
8:AI:23:GLY:O	8:AI:61:ASP:HB2	2.12	0.49
10:AK:88:PRO:HA	10:AK:92:ARG:NE	2.27	0.49
17:AR:46:THR:CG2	17:AR:51:GLN:HB2	2.42	0.49
35:BN:125:ALA:HA	47:B0:51:ARG:HH12	1.77	0.49
50:B3:24:LYS:HZ1	50:B3:29:ARG:HH12	1.59	0.49
23:BB:1051:G:H2'	23:BB:1052:C:C6	2.47	0.49
23:BB:1082:U:C4	23:BB:1086:A:N1	2.79	0.49
23:BB:2026:U:H2'	23:BB:2027:G:H8	1.78	0.49
23:BB:2240:U:O2'	23:BB:2241:A:H5'	2.12	0.49
23:BB:2502:G:H5'	23:BB:2503:A:C5'	2.36	0.49
23:BB:2514:U:H2'	23:BB:2515:C:H6	1.77	0.49
23:BB:898:C:H2'	23:BB:899:A:C8	2.46	0.49
25:BC:30:ALA:H	25:BC:31:PRO:CD	2.12	0.49
23:BB:773:U:H4'	25:BC:47:ARG:CB	2.41	0.49
26:BD:126:ASN:O	26:BD:127:PHE:CB	2.50	0.49
26:BD:7:LYS:HD3	26:BD:197:THR:H	1.76	0.49
26:BD:84:LEU:O	26:BD:84:LEU:HG	2.12	0.49
29:BG:16:VAL:HG12	29:BG:17:LYS:N	2.27	0.49
29:BG:17:LYS:HB3	29:BG:24:THR:CG2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BI:96:LYS:N	52:BI:96:LYS:HD2	2.27	0.49
32:BK:33:ALA:HB3	32:BK:39:ILE:HD11	1.94	0.49
33:BL:111:ILE:N	33:BL:111:ILE:HD13	2.26	0.49
33:BL:56:PRO:HA	33:BL:60:ARG:CZ	2.41	0.49
33:BL:82:LEU:C	33:BL:86:GLU:HG3	2.33	0.49
36:BO:87:ILE:HG22	36:BO:114:GLY:HA2	1.94	0.49
37:BP:36:LYS:HD3	37:BP:37:LYS:H	1.76	0.49
38:BQ:112:ALA:HA	38:BQ:115:ALA:HB3	1.94	0.49
40:BS:86:MET:O	40:BS:94:ASP:HB3	2.12	0.49
41:BT:66:LYS:H	41:BT:77:ARG:HA	1.77	0.49
43:BW:30:VAL:HG12	43:BW:31:LEU:N	2.27	0.49
43:BW:47:GLY:N	43:BW:67:LYS:NZ	2.60	0.49
44:BX:23:ARG:HH21	44:BX:27:ASN:CG	2.15	0.49
1:CA:1121:U:H2'	1:CA:1122:U:C6	2.47	0.49
1:CA:1285:A:O2'	1:CA:1286:U:H2'	2.12	0.49
1:CA:189:A:H2'	1:CA:190:A:C8	2.47	0.49
20:CB:130:LYS:HE2	20:CB:133:ALA:HB3	1.94	0.49
2:CC:52:SER:HA	2:CC:114:LEU:HG	1.95	0.49
9:CJ:37:ARG:C	9:CJ:37:ARG:NE	2.66	0.49
10:CK:86:LYS:HB2	10:CK:112:VAL:HG23	1.94	0.49
12:CM:15:VAL:HG13	12:CM:40:GLU:CB	2.43	0.49
12:CM:2:ARG:HB3	12:CM:6:ILE:H	1.77	0.49
49:D2:12:ARG:HG2	49:D2:46:LYS:HA	1.94	0.49
23:DB:1309:G:H4'	49:D2:7:PRO:CB	2.41	0.49
50:D3:34:LYS:N	50:D3:34:LYS:HD2	2.26	0.49
23:DB:1039:A:H2'	23:DB:1040:A:H8	1.75	0.49
23:DB:1099:G:C5'	52:DI:4:VAL:HG12	2.42	0.49
23:DB:160:A:N6	23:DB:167:A:H1'	2.26	0.49
23:DB:1747:U:H2'	23:DB:1748:C:C6	2.47	0.49
23:DB:2080:A:H2'	23:DB:2081:U:H6	1.76	0.49
23:DB:2240:U:O2'	23:DB:2241:A:H5'	2.11	0.49
23:DB:455:C:C4	23:DB:472:A:H2'	2.47	0.49
23:DB:572:A:H5''	23:DB:573:U:OP2	2.12	0.49
25:DC:220:ARG:NH1	25:DC:220:ARG:HA	2.27	0.49
26:DD:124:ARG:HB3	26:DD:124:ARG:HH11	1.77	0.49
29:DG:84:LYS:O	29:DG:132:LEU:HG	2.12	0.49
33:DL:92:LEU:HA	33:DL:96:LYS:CG	2.41	0.49
35:DN:112:TYR:O	35:DN:113:ILE:HB	2.12	0.49
35:DN:14:SER:HA	35:DN:17:ARG:NH1	2.27	0.49
23:DB:2294:G:OP1	36:DO:9:ARG:HG3	2.12	0.49
39:DR:46:GLU:HG3	39:DR:51:VAL:HG21	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:DV:76:ASP:H	24:DV:90:ASP:HB2	1.77	0.49
43:DW:42:THR:OG1	43:DW:65:LYS:HA	2.12	0.49
1:AA:1018:G:H2'	1:AA:1019:A:C8	2.47	0.49
1:AA:114:U:H2'	1:AA:115:G:C8	2.47	0.49
1:AA:123:U:H5''	1:AA:311:C:O2'	2.12	0.49
1:AA:1437:A:H2'	1:AA:1438:G:H8	1.77	0.49
1:AA:314:C:O2'	1:AA:315:A:H5'	2.12	0.49
1:AA:947:G:H2'	1:AA:948:C:H6	1.76	0.49
20:AB:101:THR:HG23	20:AB:102:ASN:N	2.28	0.49
2:AC:122:GLN:O	2:AC:127:VAL:HG22	2.12	0.49
4:AE:73:VAL:HG12	4:AE:74:ALA:N	2.27	0.49
6:AG:49:LEU:HB2	6:AG:57:GLU:HG3	1.94	0.49
7:AH:17:GLN:HE21	7:AH:62:LEU:HG	1.76	0.49
8:AI:62:LEU:N	8:AI:62:LEU:HD22	2.26	0.49
14:AO:27:GLN:O	14:AO:31:LEU:HD23	2.12	0.49
19:AT:66:ILE:HG23	19:AT:70:LYS:HD3	1.94	0.49
23:BB:1151:A:H2'	23:BB:1152:C:C6	2.47	0.49
23:BB:1681:G:H2'	23:BB:1757:A:N1	2.27	0.49
23:BB:184:C:H2'	23:BB:185:G:C8	2.48	0.49
23:BB:2247:A:H2'	23:BB:2248:C:H6	1.77	0.49
23:BB:2249:U:H4'	23:BB:2275:C:C5	2.47	0.49
23:BB:2412:A:H2'	23:BB:2413:G:O4'	2.11	0.49
23:BB:2484:G:HO2'	34:BM:45:GLN:CD	2.14	0.49
23:BB:2648:G:H2'	23:BB:2649:C:C6	2.47	0.49
23:BB:592:A:H2'	23:BB:593:U:C6	2.47	0.49
23:BB:663:G:OP1	33:BL:27:LEU:HD11	2.12	0.49
23:BB:2724:U:OP1	26:BD:116:LYS:HD2	2.12	0.49
26:BD:128:ARG:HH11	26:BD:144:GLY:HA2	1.77	0.49
26:BD:17:GLU:HA	37:BP:80:VAL:HG21	1.95	0.49
28:BF:18:GLU:O	28:BF:19:PHE:CB	2.59	0.49
28:BF:41:GLU:HG3	28:BF:49:LEU:HD12	1.94	0.49
31:BJ:103:ILE:O	31:BJ:103:ILE:HG22	2.12	0.49
31:BJ:56:VAL:HG12	31:BJ:57:LEU:H	1.77	0.49
31:BJ:72:LYS:O	31:BJ:74:TYR:N	2.46	0.49
33:BL:123:ARG:HH22	33:BL:141:LYS:NZ	2.10	0.49
34:BM:89:VAL:HG22	34:BM:89:VAL:O	2.12	0.49
26:BD:25:THR:HB	37:BP:5:LYS:HZ3	1.77	0.49
39:BR:65:ALA:O	39:BR:66:HIS:HB2	2.12	0.49
39:BR:85:LYS:O	39:BR:85:LYS:HG3	2.12	0.49
41:BT:69:ARG:N	41:BT:75:GLY:HA3	2.28	0.49
24:BV:26:PHE:CE2	24:BV:44:HIS:HA	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:BT:8:LEU:CB	44:BX:19:LEU:HD11	2.32	0.49
1:CA:1229:A:H2'	1:CA:1230:C:C6	2.47	0.49
1:CA:1343:G:H2'	1:CA:1344:C:H6	1.78	0.49
1:CA:418:C:H2'	1:CA:419:C:H6	1.78	0.49
1:CA:626:G:H2'	1:CA:627:G:C8	2.48	0.49
2:CC:88:LYS:O	2:CC:88:LYS:HE3	2.13	0.49
12:CM:79:LEU:HD22	12:CM:86:ARG:NE	2.27	0.49
12:CM:95:PRO:HB2	12:CM:99:GLN:OE1	2.11	0.49
13:CN:61:ASN:HB2	13:CN:72:PHE:CZ	2.47	0.49
15:CP:22:ALA:HA	15:CP:33:ILE:HG13	1.95	0.49
15:CP:78:VAL:O	15:CP:78:VAL:HG13	2.12	0.49
16:CQ:13:SER:O	16:CQ:20:ILE:HG13	2.13	0.49
16:CQ:17:GLU:C	16:CQ:19:SER:H	2.15	0.49
16:CQ:56:ASP:HA	16:CQ:81:ALA:CA	2.42	0.49
47:D0:12:ARG:NH2	47:D0:16:ARG:HG3	2.27	0.49
23:DB:1038:G:H2'	23:DB:1039:A:H8	1.76	0.49
23:DB:1152:C:O2'	23:DB:1153:C:H5'	2.11	0.49
23:DB:1299:G:C4'	23:DB:1301:A:H1'	2.40	0.49
23:DB:2244:U:H1'	23:DB:2434:A:C4	2.48	0.49
23:DB:2776:A:H4'	23:DB:2777:G:O5'	2.13	0.49
23:DB:2803:G:H2'	23:DB:2804:U:C6	2.47	0.49
23:DB:2876:G:H4'	37:DP:2:ASN:HB3	1.93	0.49
23:DB:335:C:O2'	23:DB:336:C:H5'	2.12	0.49
23:DB:521:U:H2'	23:DB:522:A:H8	1.76	0.49
23:DB:764:A:H5''	25:DC:208:GLY:CA	2.42	0.49
26:DD:148:GLN:HB3	26:DD:151:THR:HG23	1.95	0.49
27:DE:127:GLU:CA	27:DE:157:LEU:HD13	2.32	0.49
28:DF:125:GLY:O	28:DF:157:THR:HB	2.13	0.49
30:DH:54:LEU:O	30:DH:58:LEU:HD23	2.12	0.49
52:DI:19:PRO:HB2	52:DI:22:PRO:HD2	1.93	0.49
52:DI:63:ASP:OD1	52:DI:65:SER:HB2	2.12	0.49
31:DJ:90:GLU:O	31:DJ:93:ILE:HG13	2.12	0.49
32:DK:99:ILE:HD12	32:DK:118:LEU:HD12	1.94	0.49
33:DL:108:ALA:HB1	33:DL:125:LEU:HD13	1.94	0.49
34:DM:119:LEU:HD22	34:DM:119:LEU:H	1.78	0.49
34:DM:20:LEU:HD22	34:DM:38:ARG:HG2	1.94	0.49
34:DM:71:LYS:O	34:DM:73:ILE:N	2.45	0.49
35:DN:5:LYS:HA	35:DN:5:LYS:CE	2.34	0.49
40:DS:24:ILE:HG23	40:DS:35:ILE:HG21	1.93	0.49
42:DU:66:VAL:HG13	42:DU:67:SER:N	2.27	0.49
45:DY:23:LEU:HD13	45:DY:28:LEU:HD23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:DY:43:ILE:HD11	45:DY:47:ILE:HD11	1.94	0.49
23:DB:398:C:OP1	46:DZ:49:ARG:NH1	2.45	0.49
1:AA:182:A:H1'	1:AA:183:C:C5	2.48	0.49
1:AA:205:A:H2'	1:AA:206:C:C6	2.44	0.49
1:AA:317:U:H2'	1:AA:318:G:C8	2.47	0.49
1:AA:366:A:O2'	1:AA:367:U:P	2.70	0.49
1:AA:576:C:OP2	1:AA:576:C:H3'	2.12	0.49
1:AA:662:U:O2'	1:AA:836:G:H5''	2.12	0.49
2:AC:143:LEU:HD22	2:AC:143:LEU:H	1.76	0.49
5:AF:39:LEU:HD13	5:AF:40:GLU:H	1.77	0.49
7:AH:39:LEU:HB2	7:AH:45:ILE:CD1	2.43	0.49
8:AI:32:ARG:HD3	8:AI:37:TYR:HD1	1.75	0.49
1:AA:1048:G:H4'	13:AN:2:LYS:HZ1	1.76	0.49
15:AP:29:ASN:HD22	15:AP:29:ASN:N	2.11	0.49
18:AS:35:ARG:HG2	18:AS:50:VAL:HG13	1.94	0.49
47:B0:29:VAL:HG13	47:B0:30:ASP:N	2.26	0.49
22:BA:74:U:H2'	22:BA:75:G:O4'	2.12	0.49
23:BB:1169:A:O2'	23:BB:1170:C:H5'	2.12	0.49
23:BB:1684:G:H2'	23:BB:1685:C:H6	1.77	0.49
23:BB:2528:U:O2'	23:BB:2529:G:H3'	2.13	0.49
23:BB:2590:A:H2'	23:BB:2591:C:H6	1.76	0.49
23:BB:650:C:O3'	50:B3:16:THR:HB	2.12	0.49
23:BB:836:G:H2'	23:BB:837:C:H6	1.77	0.49
25:BC:209:ALA:HA	25:BC:213:ARG:CZ	2.43	0.49
26:BD:4:LEU:HD23	26:BD:4:LEU:H	1.77	0.49
23:BB:38:A:H1'	27:BE:44:ARG:HA	1.93	0.49
52:BI:116:MET:SD	52:BI:124:MET:HB2	2.52	0.49
31:BJ:15:TRP:HB3	31:BJ:53:TYR:CD2	2.48	0.49
38:BQ:77:LYS:HB3	38:BQ:116:LEU:HD11	1.93	0.49
38:BQ:27:ARG:N	38:BQ:32:ARG:HH21	2.11	0.49
40:BS:29:VAL:HG21	40:BS:69:LEU:C	2.33	0.49
41:BT:16:VAL:HG22	41:BT:16:VAL:O	2.13	0.49
46:BZ:5:ILE:HG23	46:BZ:63:ARG:NE	2.27	0.49
1:CA:1031:C:H5'	1:CA:1032:G:N9	2.28	0.49
1:CA:1121:U:O2'	1:CA:1122:U:H5'	2.13	0.49
1:CA:1237:C:OP1	1:CA:1238:A:H1'	2.11	0.49
1:CA:1243:C:H2'	1:CA:1244:G:C8	2.48	0.49
1:CA:1526:G:O5'	21:CU:38:GLU:HB2	2.12	0.49
1:CA:272:C:H2'	1:CA:273:U:C6	2.47	0.49
1:CA:712:A:O2'	1:CA:713:G:H5'	2.12	0.49
20:CB:67:LEU:HD22	20:CB:157:PRO:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:8:A:C5	3:CD:205:LYS:HA	2.47	0.49
4:CE:114:LEU:HD13	4:CE:122:VAL:HG21	1.94	0.49
4:CE:81:GLN:NE2	4:CE:148:SER:HA	2.26	0.49
5:CF:18:VAL:O	5:CF:22:ILE:HG13	2.13	0.49
5:CF:85:ILE:HG22	5:CF:86:ARG:N	2.27	0.49
7:CH:10:LEU:HD11	7:CH:126:CYS:SG	2.52	0.49
12:CM:80:MET:HA	12:CM:87:GLY:HA3	1.94	0.49
14:CO:25:GLU:OE2	14:CO:76:ARG:HD3	2.12	0.49
16:CQ:58:VAL:HB	16:CQ:74:LEU:HD23	1.94	0.49
48:D1:34:GLU:HG2	48:D1:50:GLU:OE1	2.11	0.49
22:DA:63:C:H2'	22:DA:64:G:H8	1.77	0.49
23:DB:2047:C:O2'	23:DB:2048:G:H5'	2.12	0.49
23:DB:2314:A:H2'	23:DB:2315:G:H8	1.77	0.49
23:DB:2425:A:O2'	23:DB:2426:A:OP2	2.28	0.49
23:DB:244:A:H2'	23:DB:245:G:O4'	2.12	0.49
23:DB:2654:A:N1	23:DB:2665:A:H5''	2.27	0.49
23:DB:2743:U:H2'	23:DB:2744:G:C5'	2.40	0.49
23:DB:2860:A:H2'	23:DB:2861:U:O4'	2.12	0.49
23:DB:41:C:O2'	23:DB:42:A:H5'	2.13	0.49
23:DB:592:A:H2'	23:DB:593:U:C6	2.47	0.49
25:DC:37:SER:HB2	25:DC:62:ARG:HG3	1.94	0.49
27:DE:191:ASP:HA	27:DE:194:LYS:NZ	2.27	0.49
29:DG:17:LYS:O	29:DG:17:LYS:HE3	2.12	0.49
33:DL:18:ARG:NH1	33:DL:21:ARG:HD3	2.26	0.49
23:DB:2393:U:H4'	33:DL:60:ARG:O	2.12	0.49
33:DL:62:PRO:HG3	50:D3:12:ARG:CZ	2.43	0.49
34:DM:5:LYS:HG2	34:DM:68:PHE:HE1	1.78	0.49
23:DB:1199:U:O2'	38:DQ:2:ARG:CB	2.60	0.49
23:DB:536:G:C5'	38:DQ:52:ARG:NH2	2.75	0.49
39:DR:37:GLU:HG2	39:DR:62:GLU:C	2.33	0.49
39:DR:69:GLY:HA2	39:DR:96:VAL:HG13	1.94	0.49
41:DT:67:VAL:HG12	41:DT:68:LYS:N	2.26	0.49
44:DX:18:LEU:O	44:DX:22:LEU:HB2	2.13	0.49
44:DX:22:LEU:HB3	44:DX:25:GLN:OE1	2.12	0.49
1:AA:1221:G:O3'	18:AS:76:THR:HG21	2.12	0.49
1:AA:1386:G:O2'	1:AA:1387:G:H5'	2.12	0.49
1:AA:997:U:H2'	1:AA:998:C:C6	2.46	0.49
20:AB:119:GLN:HE22	20:AB:127:LYS:HD3	1.77	0.49
4:AE:87:VAL:CG1	4:AE:92:ARG:HD3	2.42	0.49
10:AK:85:VAL:O	10:AK:111:ASP:HA	2.12	0.49
12:AM:53:ASP:HA	12:AM:56:ARG:HD2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AO:66:LEU:HB3	14:AO:77:TYR:HE1	1.77	0.49
14:AO:68:TYR:CE1	14:AO:72:LYS:HE3	2.47	0.49
23:BB:1255:U:O2'	27:BE:67:ARG:HB3	2.12	0.49
23:BB:1501:G:O2'	23:BB:1502:A:H5'	2.12	0.49
23:BB:2082:A:O2'	23:BB:2083:G:H5'	2.13	0.49
23:BB:2249:U:H4'	23:BB:2275:C:H5	1.76	0.49
23:BB:2314:A:H2'	23:BB:2315:G:C8	2.47	0.49
23:BB:2831:G:O4'	23:BB:2883:A:C2	2.65	0.49
23:BB:69:C:H2'	23:BB:70:G:C8	2.47	0.49
25:BC:66:PHE:HE2	25:BC:104:LEU:HD21	1.77	0.49
25:BC:12:ARG:HB3	25:BC:18:VAL:HB	1.94	0.49
25:BC:204:LEU:CD1	25:BC:211:ARG:HH11	2.25	0.49
26:BD:157:LYS:HE2	26:BD:157:LYS:HA	1.94	0.49
27:BE:154:ASP:OD1	27:BE:156:ASN:HB2	2.13	0.49
28:BF:90:LEU:O	28:BF:95:MET:HG3	2.12	0.49
29:BG:6:ALA:HB3	29:BG:7:PRO:HD3	1.93	0.49
31:BJ:17:VAL:HG22	31:BJ:55:ILE:CG2	2.42	0.49
34:BM:62:LYS:CB	34:BM:104:GLU:HB2	2.42	0.49
34:BM:67:VAL:HG21	34:BM:95:LEU:HD11	1.94	0.49
23:BB:2000:C:OP1	35:BN:2:ARG:NE	2.45	0.49
36:BO:8:ILE:CG2	36:BO:9:ARG:N	2.75	0.49
42:BU:3:LYS:HZ3	42:BU:3:LYS:HB2	1.72	0.49
43:BW:16:GLU:HG3	43:BW:37:VAL:HG22	1.93	0.49
43:BW:43:LYS:HG2	43:BW:76:ARG:NE	2.25	0.49
44:BX:19:LEU:HD12	44:BX:19:LEU:H	1.77	0.49
1:CA:1162:C:H2'	1:CA:1163:A:C8	2.47	0.49
1:CA:1163:A:H2'	1:CA:1164:G:C8	2.46	0.49
1:CA:1054:C:H1'	1:CA:1196:A:C5	2.48	0.49
1:CA:1238:A:H2	1:CA:1241:G:N3	2.10	0.49
1:CA:219:U:H2'	1:CA:220:G:H8	1.77	0.49
1:CA:33:A:H2'	1:CA:34:C:H6	1.77	0.49
3:CD:106:PHE:HD1	3:CD:158:LEU:HD21	1.78	0.49
3:CD:96:ARG:O	3:CD:99:ASN:HB3	2.12	0.49
4:CE:88:HIS:CE1	4:CE:137:ARG:HG2	2.48	0.49
5:CF:42:TRP:NE1	5:CF:61:LEU:HD23	2.27	0.49
9:CJ:102:LEU:N	9:CJ:102:LEU:HD22	2.27	0.49
12:CM:22:TYR:HB3	12:CM:69:ARG:NH2	2.27	0.49
12:CM:78:ARG:HG3	28:DF:111:ARG:NH1	2.26	0.49
10:CK:112:VAL:HA	17:CR:72:ARG:HD3	1.94	0.49
47:D0:11:LYS:O	47:D0:15:ARG:HB2	2.13	0.49
48:D1:29:LYS:CB	48:D1:30:PRO:HD3	2.17	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:D4:3:VAL:CG1	51:D4:4:ARG:H	2.13	0.49
23:DB:1135:C:N4	23:DB:1139:G:C6	2.81	0.49
23:DB:15:G:O2'	23:DB:16:C:H5'	2.13	0.49
23:DB:21:A:H2'	23:DB:22:C:C6	2.48	0.49
23:DB:2377:A:C2	36:DO:92:PHE:HE1	2.30	0.49
23:DB:2528:U:O2'	23:DB:2529:G:H3'	2.13	0.49
23:DB:378:C:O2'	23:DB:379:G:H5'	2.12	0.49
23:DB:587:C:H4'	23:DB:588:U:C6	2.48	0.49
23:DB:721:A:H2'	23:DB:722:A:C8	2.48	0.49
23:DB:827:U:H5'	23:DB:828:U:O5'	2.12	0.49
25:DC:86:ARG:HG3	25:DC:86:ARG:O	2.12	0.49
27:DE:46:GLN:HG3	27:DE:49:ARG:NH2	2.28	0.49
28:DF:40:GLY:H	28:DF:84:ILE:CG2	2.26	0.49
30:DH:125:THR:HA	30:DH:146:VAL:CB	2.32	0.49
32:DK:64:ARG:O	32:DK:65:THR:HG23	2.13	0.49
33:DL:77:ILE:CG2	33:DL:92:LEU:HG	2.42	0.49
41:DT:92:ASN:C	41:DT:93:LEU:HD12	2.32	0.49
43:DW:42:THR:HG22	43:DW:67:LYS:O	2.13	0.49
1:AA:1492:A:H2'	1:AA:1493:A:O4'	2.12	0.49
1:AA:301:G:H2'	1:AA:302:G:C8	2.48	0.49
1:AA:35:G:H2'	1:AA:36:C:H6	1.78	0.49
1:AA:382:A:H2'	1:AA:383:A:C8	2.48	0.49
1:AA:724:G:O2'	1:AA:725:G:H5'	2.11	0.49
20:AB:173:LYS:HD3	20:AB:173:LYS:C	2.32	0.49
8:AI:24:ASN:O	8:AI:61:ASP:HA	2.12	0.49
10:AK:69:CYS:O	10:AK:73:VAL:HG23	2.12	0.49
11:AL:34:THR:HG21	11:AL:53:ARG:CZ	2.43	0.49
16:AQ:19:SER:O	16:AQ:20:ILE:HG23	2.12	0.49
23:BB:1292:G:H2'	23:BB:1293:C:C6	2.48	0.49
23:BB:1322:A:C2'	23:BB:1323:C:H5'	2.43	0.49
23:BB:139:U:H5''	23:BB:140:C:C4	2.47	0.49
23:BB:142:A:H1'	41:BT:2:ILE:HG21	1.93	0.49
23:BB:1562:U:H2'	23:BB:1563:U:C6	2.47	0.49
23:BB:1727:C:H2'	23:BB:1728:C:H6	1.74	0.49
23:BB:1881:C:H2'	23:BB:1882:U:O4'	2.13	0.49
23:BB:1948:G:O2'	23:BB:1949:G:H5'	2.12	0.49
23:BB:2287:A:H4'	23:BB:2287:A:OP1	2.10	0.49
25:BC:70:LYS:HG2	25:BC:73:ILE:HD13	1.94	0.49
26:BD:15:PHE:CD2	26:BD:18:ASP:HB2	2.47	0.49
26:BD:170:VAL:HG11	26:BD:194:PRO:HB2	1.94	0.49
26:BD:83:ARG:HG2	26:BD:84:LEU:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:21:TYR:OH	28:BF:28:PRO:HD3	2.13	0.49
28:BF:41:GLU:OE2	28:BF:48:LEU:HB3	2.13	0.49
29:BG:40:VAL:CB	29:BG:63:GLN:HE22	2.25	0.49
30:BH:80:ILE:HD12	30:BH:102:ALA:CB	2.43	0.49
23:BB:1190:G:OP1	33:BL:41:ARG:CZ	2.60	0.49
33:BL:57:LEU:C	33:BL:59:ARG:H	2.13	0.49
33:BL:60:ARG:C	50:B3:11:LYS:HZ3	2.16	0.49
35:BN:102:PHE:CD2	35:BN:102:PHE:N	2.80	0.49
37:BP:13:LYS:HA	37:BP:15:ASP:OD1	2.13	0.49
38:BQ:63:ARG:CB	38:BQ:63:ARG:HH11	2.26	0.49
44:BX:18:LEU:HD12	44:BX:19:LEU:HD12	1.94	0.49
44:BX:22:LEU:HD22	44:BX:26:PHE:N	2.23	0.49
44:BX:30:MET:CG	44:BX:31:GLN:H	2.24	0.49
44:BX:46:VAL:HG13	44:BX:47:ARG:N	2.28	0.49
46:BZ:59:ARG:HB2	46:BZ:61:ASN:H	1.78	0.49
1:CA:1228:C:OP1	12:CM:112:ARG:HA	2.13	0.49
1:CA:147:G:H2'	1:CA:148:G:C8	2.48	0.49
1:CA:366:A:O2'	1:CA:367:U:P	2.70	0.49
1:CA:662:U:O2'	1:CA:836:G:H5''	2.12	0.49
1:CA:693:G:H2'	1:CA:694:A:O4'	2.12	0.49
2:CC:64:ARG:HA	2:CC:99:GLN:HB2	1.93	0.49
3:CD:130:ASN:N	3:CD:130:ASN:HD22	2.06	0.49
9:CJ:87:LEU:N	9:CJ:87:LEU:HD22	2.27	0.49
10:CK:31:VAL:HG21	10:CK:66:ALA:HA	1.94	0.49
18:CS:43:MET:C	18:CS:46:LEU:HD23	2.32	0.49
22:DA:116:G:H4'	36:DO:54:VAL:CG2	2.42	0.49
23:DB:1082:U:H2'	23:DB:1083:U:O4'	2.12	0.49
23:DB:1214:A:H2'	23:DB:1215:G:O4'	2.12	0.49
23:DB:1256:G:H1'	27:DE:77:ILE:HD11	1.95	0.49
23:DB:1292:G:H2'	23:DB:1293:C:C6	2.48	0.49
23:DB:185:G:H4'	23:DB:218:A:H4'	1.94	0.49
23:DB:1881:C:H2'	23:DB:1882:U:O4'	2.12	0.49
23:DB:1909:C:H2'	23:DB:1910:G:H8	1.77	0.49
23:DB:2518:A:H2'	23:DB:2518:A:N3	2.26	0.49
23:DB:2590:A:H2'	23:DB:2591:C:H6	1.76	0.49
23:DB:2648:G:H2'	23:DB:2649:C:C6	2.47	0.49
23:DB:2717:C:O2'	37:DP:95:LYS:HE3	2.13	0.49
23:DB:303:G:H2'	23:DB:304:U:C6	2.47	0.49
23:DB:365:U:H2'	23:DB:366:C:O4'	2.12	0.49
23:DB:465:G:H4'	49:D2:12:ARG:NH2	2.27	0.49
25:DC:143:VAL:O	25:DC:152:GLN:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:323:C:O2'	27:DE:163:ASN:C	2.50	0.49
27:DE:169:VAL:HG22	27:DE:170:ARG:N	2.28	0.49
27:DE:20:GLY:HA2	27:DE:106:LYS:HD2	1.95	0.49
23:DB:2093:G:OP2	30:DH:23:ALA:HB3	2.12	0.49
30:DH:81:ALA:HA	30:DH:147:VAL:O	2.12	0.49
52:DI:7:TYR:HA	52:DI:59:THR:HA	1.95	0.49
23:DB:1995:U:H1'	32:DK:3:GLN:HE22	1.78	0.49
33:DL:108:ALA:O	33:DL:109:LYS:CB	2.61	0.49
34:DM:11:LYS:O	34:DM:12:MET:HB2	2.13	0.49
34:DM:42:THR:HG22	34:DM:45:GLN:NE2	2.27	0.49
35:DN:35:LYS:HD3	35:DN:112:TYR:CE1	2.47	0.49
38:DQ:89:ILE:HG22	38:DQ:90:ASP:N	2.26	0.49
39:DR:11:GLN:C	39:DR:21:ARG:HH22	2.14	0.49
41:DT:26:LYS:O	41:DT:27:SER:HB3	2.13	0.49
42:DU:10:VAL:HB	42:DU:68:ASN:ND2	2.27	0.49
24:DV:70:ILE:CD1	24:DV:71:LYS:H	2.23	0.49
23:DB:2386:A:H4'	43:DW:38:ARG:HD3	1.94	0.49
46:DZ:59:ARG:HA	46:DZ:62:LYS:HB2	1.94	0.49
1:AA:693:G:H2'	1:AA:694:A:O4'	2.12	0.49
1:AA:853:C:O2'	1:AA:854:U:H5'	2.12	0.49
20:AB:132:GLU:HG3	20:AB:136:ARG:HD2	1.94	0.49
20:AB:139:GLU:HG2	20:AB:143:LEU:CD1	2.42	0.49
20:AB:57:ASN:HB2	20:AB:219:THR:HB	1.95	0.49
2:AC:110:LEU:HB3	2:AC:203:LYS:HZ3	1.77	0.49
2:AC:11:LEU:HD13	2:AC:17:TRP:HE1	1.76	0.49
8:AI:96:GLU:O	8:AI:99:LYS:HB3	2.12	0.49
21:AU:20:ARG:O	21:AU:20:ARG:HG2	2.13	0.49
23:BB:1685:C:H2'	23:BB:1686:C:H6	1.78	0.49
23:BB:1853:A:N1	23:BB:2087:G:H1'	2.27	0.49
23:BB:209:C:O2'	23:BB:210:C:H5'	2.12	0.49
23:BB:2654:A:N1	23:BB:2665:A:H5''	2.27	0.49
23:BB:2840:C:H2'	23:BB:2841:C:C6	2.48	0.49
23:BB:600:G:H1'	27:BE:100:MET:HG3	1.95	0.49
23:BB:991:C:H2'	23:BB:992:C:C6	2.47	0.49
26:BD:130:GLN:HE21	26:BD:138:LEU:HD12	1.71	0.49
23:BB:2786:U:C4'	26:BD:66:GLY:HA2	2.43	0.49
26:BD:34:VAL:CA	26:BD:89:GLU:HB2	2.34	0.49
23:BB:1205:A:O2'	27:BE:164:LEU:HD21	2.13	0.49
29:BG:139:VAL:O	29:BG:143:VAL:HG23	2.13	0.49
30:BH:67:ALA:O	30:BH:71:LYS:HG3	2.13	0.49
30:BH:89:LYS:HZ3	30:BH:90:LEU:H	1.61	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BI:17:ALA:O	52:BI:18:ASN:HB3	2.12	0.49
52:BI:63:ASP:C	52:BI:65:SER:H	2.15	0.49
31:BJ:128:ASN:ND2	31:BJ:129:GLU:H	2.11	0.49
35:BN:34:ILE:HD13	35:BN:34:ILE:H	1.76	0.49
37:BP:23:ASP:O	37:BP:25:VAL:N	2.46	0.49
39:BR:69:GLY:HA2	39:BR:96:VAL:HA	1.94	0.49
23:BB:25:U:OP1	40:BS:8:ARG:NH2	2.45	0.49
1:CA:1320:C:H41	18:CS:36:ARG:CD	2.25	0.49
1:CA:1324:A:H2'	1:CA:1325:C:C6	2.48	0.49
1:CA:317:U:H2'	1:CA:318:G:C8	2.47	0.49
1:CA:966:G:H2'	1:CA:967:C:C6	2.47	0.49
4:CE:85:LYS:HE3	4:CE:94:PHE:HB2	1.95	0.49
5:CF:84:VAL:HG22	5:CF:85:ILE:N	2.28	0.49
10:CK:31:VAL:HG21	10:CK:66:ALA:CA	2.42	0.49
12:CM:102:LYS:HZ2	12:CM:102:LYS:HB2	1.77	0.49
16:CQ:46:HIS:CG	16:CQ:47:ASP:N	2.81	0.49
48:D1:7:LYS:HG2	48:D1:26:LYS:HB3	1.94	0.49
23:DB:1259:G:H2'	23:DB:1260:A:C8	2.48	0.49
23:DB:1684:G:H2'	23:DB:1685:C:H6	1.77	0.49
23:DB:2230:G:H5''	46:DZ:28:VAL:HG21	1.94	0.49
23:DB:2734:A:C2'	23:DB:2735:G:H5'	2.43	0.49
23:DB:345:A:H1'	23:DB:346:A:H2	1.77	0.49
23:DB:383:C:N4	23:DB:385:C:H2'	2.27	0.49
23:DB:778:G:H5''	25:DC:48:ILE:HG21	1.95	0.49
23:DB:845:A:C2	23:DB:847:U:H1'	2.47	0.49
25:DC:163:ILE:O	25:DC:164:VAL:HG13	2.13	0.49
25:DC:169:ALA:O	25:DC:170:TYR:HB2	2.13	0.49
25:DC:231:HIS:CE1	25:DC:242:HIS:HD1	2.31	0.49
26:DD:205:PRO:HG2	26:DD:206:ALA:H	1.77	0.49
27:DE:115:GLN:OE1	27:DE:184:ASP:HB2	2.13	0.49
29:DG:79:THR:HG23	29:DG:80:GLU:H	1.78	0.49
30:DH:82:SER:C	30:DH:83:LYS:HD2	2.33	0.49
30:DH:86:ASP:C	30:DH:88:GLY:H	2.16	0.49
52:DI:108:ILE:CG2	52:DI:128:ILE:HD13	2.43	0.49
31:DJ:7:LYS:HE3	31:DJ:47:HIS:HD2	1.76	0.49
31:DJ:58:ASN:C	31:DJ:60:ASP:H	2.16	0.49
32:DK:63:VAL:HG11	32:DK:103:VAL:HG12	1.95	0.49
26:DD:17:GLU:CG	37:DP:80:VAL:HB	2.42	0.49
37:DP:79:VAL:O	37:DP:81:ASP:N	2.46	0.49
40:DS:2:GLU:CB	40:DS:108:SER:HA	2.41	0.49
42:DU:8:ASP:O	42:DU:10:VAL:HG13	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DW:10:ARG:HA	43:DW:10:ARG:HH11	1.77	0.49
1:AA:157:U:O2'	1:AA:158:G:H5'	2.13	0.49
1:AA:218:U:H2'	1:AA:219:U:H6	1.78	0.49
1:AA:272:C:H2'	1:AA:273:U:C6	2.48	0.49
1:AA:627:G:H2'	1:AA:628:G:H8	1.78	0.49
1:AA:692:U:H2'	1:AA:694:A:OP2	2.13	0.49
1:AA:846:G:H2'	1:AA:846:G:N3	2.27	0.49
1:AA:975:A:H4'	1:AA:976:G:OP2	2.13	0.49
20:AB:27:LYS:H	20:AB:28:PRO:CD	2.25	0.49
20:AB:67:LEU:HA	20:AB:89:PHE:O	2.11	0.49
4:AE:104:ILE:O	4:AE:104:ILE:HG23	2.13	0.49
4:AE:45:VAL:CG1	4:AE:116:VAL:HG23	2.43	0.49
4:AE:45:VAL:HG13	4:AE:117:ALA:HA	1.95	0.49
4:AE:67:ARG:O	4:AE:70:MET:HG3	2.13	0.49
5:AF:1:MET:SD	5:AF:67:PRO:HD3	2.53	0.49
9:AJ:53:ILE:HB	9:AJ:61:ALA:HB1	1.94	0.49
13:AN:24:ALA:HB1	13:AN:27:LYS:HZ1	1.78	0.49
16:AQ:10:ARG:NH1	16:AQ:11:VAL:HB	2.28	0.49
16:AQ:20:ILE:CG1	16:AQ:45:VAL:HB	2.37	0.49
16:AQ:60:ILE:HB	16:AQ:72:TRP:HE3	1.78	0.49
50:B3:12:ARG:CB	50:B3:23:HIS:HA	2.34	0.49
51:B4:23:ILE:HD12	51:B4:35:GLN:HG2	1.95	0.49
23:BB:1533:C:C2'	23:BB:1534:U:H5'	2.42	0.49
23:BB:1568:G:C5'	25:BC:60:ALA:HB3	2.42	0.49
23:BB:1581:G:O2'	23:BB:1582:C:H5'	2.13	0.49
23:BB:1821:A:H2'	23:BB:1822:C:H6	1.76	0.49
23:BB:1869:G:HO2'	23:BB:1871:A:H62	1.60	0.49
23:BB:2080:A:H2'	23:BB:2081:U:C6	2.48	0.49
23:BB:2133:G:N3	23:BB:2133:G:C2'	2.73	0.49
23:BB:2336:A:C1'	23:BB:2337:G:OP1	2.56	0.49
23:BB:2500:U:H5'	23:BB:2501:C:OP2	2.12	0.49
23:BB:354:A:C2'	23:BB:355:U:H5''	2.42	0.49
23:BB:638:G:H2'	23:BB:639:U:C6	2.48	0.49
25:BC:70:LYS:HG2	25:BC:71:ASP:N	2.28	0.49
26:BD:117:GLY:O	26:BD:165:MET:HB3	2.13	0.49
26:BD:29:VAL:O	26:BD:51:THR:HG21	2.13	0.49
27:BE:152:GLU:O	27:BE:153:LEU:HB2	2.12	0.49
23:BB:673:C:H5''	27:BE:76:PRO:HD2	1.94	0.49
28:BF:165:GLY:O	28:BF:167:ALA:N	2.44	0.49
28:BF:2:LYS:HA	28:BF:2:LYS:HE2	1.92	0.49
31:BJ:118:MET:O	31:BJ:121:LYS:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BR:40:MET:CA	39:BR:54:VAL:HG11	2.43	0.49
39:BR:97:LYS:O	39:BR:98:ILE:HB	2.13	0.49
43:BW:24:ARG:HA	43:BW:57:THR:HG23	1.93	0.49
1:CA:1014:A:H2	1:CA:1219:A:H1'	1.76	0.49
1:CA:1281:C:H5'	1:CA:1282:C:C5	2.43	0.49
1:CA:1312:G:H2'	1:CA:1313:U:C6	2.48	0.49
1:CA:1366:C:H2'	1:CA:1367:C:C6	2.47	0.49
1:CA:1522:U:O2'	1:CA:1523:G:H5'	2.13	0.49
1:CA:376:G:C4'	15:CP:5:ARG:HD3	2.42	0.49
1:CA:708:C:H2'	1:CA:709:U:C6	2.48	0.49
20:CB:14:HIS:CG	20:CB:15:PHE:N	2.80	0.49
20:CB:26:MET:HG3	20:CB:188:THR:O	2.13	0.49
4:CE:95:MET:CE	4:CE:143:LEU:HD13	2.43	0.49
7:CH:38:VAL:HG23	7:CH:111:THR:HG22	1.94	0.49
10:CK:18:GLY:HA2	10:CK:35:ASP:HA	1.95	0.49
11:CL:88:ASP:C	11:CL:89:LEU:HD22	2.32	0.49
16:CQ:12:VAL:HG11	16:CQ:42:LYS:NZ	2.27	0.49
18:CS:62:THR:HG22	18:CS:65:MET:CE	2.43	0.49
19:CT:68:LYS:CA	19:CT:68:LYS:HE2	2.37	0.49
23:DB:1098:A:O2'	52:DI:4:VAL:C	2.50	0.49
23:DB:1266:G:OP1	47:D0:15:ARG:NH2	2.45	0.49
23:DB:1315:C:H2'	23:DB:1316:U:H6	1.77	0.49
23:DB:1383:A:H2	23:DB:1405:U:O2	1.96	0.49
23:DB:1904:G:H1'	23:DB:1927:A:N1	2.28	0.49
23:DB:2034:U:H5''	55:DB:3318:HOH:O	2.13	0.49
23:DB:2278:A:H62	43:DW:10:ARG:CB	2.26	0.49
23:DB:2468:A:H2'	23:DB:2476:A:C6	2.48	0.49
23:DB:2740:A:H2'	23:DB:2741:A:C8	2.47	0.49
23:DB:281:C:O2	23:DB:281:C:H2'	2.11	0.49
23:DB:2896:C:H2'	23:DB:2897:U:C6	2.47	0.49
26:DD:4:LEU:HD11	26:DD:79:LEU:HD22	1.93	0.49
29:DG:156:TYR:HA	29:DG:171:LYS:CG	2.43	0.49
31:DJ:28:LEU:C	31:DJ:28:LEU:HD13	2.32	0.49
31:DJ:64:VAL:CG1	31:DJ:65:THR:H	2.19	0.49
32:DK:70:ARG:HD3	32:DK:76:VAL:CG2	2.36	0.49
39:DR:68:ARG:HB3	39:DR:70:GLU:HG2	1.93	0.49
43:DW:4:LYS:HB2	43:DW:4:LYS:HZ3	1.78	0.49
45:DY:2:LYS:CE	45:DY:27:GLY:H	2.26	0.49
1:AA:1103:C:C5'	20:AB:96:LEU:HD12	2.42	0.49
1:AA:1121:U:O2'	1:AA:1122:U:H5'	2.13	0.49
1:AA:1371:G:OP1	8:AI:12:LYS:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:176:C:H2'	1:AA:177:G:N3	2.26	0.49
1:AA:626:G:H2'	1:AA:627:G:C8	2.48	0.49
20:AB:162:VAL:HG21	20:AB:168:GLU:HB2	1.94	0.49
2:AC:61:LYS:O	2:AC:97:PRO:HD2	2.13	0.49
1:AA:16:A:H4'	4:AE:21:SER:H	1.77	0.49
9:AJ:41:PRO:O	9:AJ:42:LEU:HB2	2.13	0.49
15:AP:43:ALA:O	15:AP:46:LYS:HD3	2.12	0.49
22:BA:82:U:O2'	22:BA:83:G:H5'	2.13	0.49
23:BB:1206:G:H2'	23:BB:1207:C:H6	1.78	0.49
23:BB:1268:A:H2'	23:BB:1269:A:O4'	2.13	0.49
23:BB:154:U:H2'	23:BB:155:A:C8	2.48	0.49
23:BB:1904:G:H1'	23:BB:1927:A:N1	2.28	0.49
23:BB:2390:U:O2'	23:BB:2391:G:H5''	2.12	0.49
23:BB:2488:G:O2'	23:BB:2489:U:H5'	2.12	0.49
23:BB:2563:U:H2'	23:BB:2565:A:OP2	2.13	0.49
23:BB:2896:C:H2'	23:BB:2897:U:H6	1.78	0.49
23:BB:967:U:H2'	23:BB:968:C:H6	1.73	0.49
25:BC:212:TRP:CZ3	25:BC:217:PRO:HD3	2.48	0.49
27:BE:29:HIS:O	27:BE:33:VAL:HG23	2.12	0.49
28:BF:55:ASP:OD1	28:BF:148:VAL:HG11	2.13	0.49
32:BK:54:LYS:HD3	32:BK:54:LYS:N	2.28	0.49
34:BM:130:PHE:HD2	34:BM:131:VAL:HG22	1.77	0.49
34:BM:71:LYS:HD3	34:BM:92:TRP:O	2.12	0.49
35:BN:13:ASN:O	35:BN:17:ARG:N	2.44	0.49
35:BN:2:ARG:O	35:BN:3:HIS:CG	2.66	0.49
35:BN:57:THR:OG1	35:BN:61:ALA:HB3	2.12	0.49
37:BP:92:ARG:HH22	37:BP:110:LYS:HE2	1.75	0.49
44:BX:8:GLU:O	44:BX:10:SER:N	2.46	0.49
1:CA:193:C:H2'	1:CA:194:C:C6	2.47	0.49
1:CA:382:A:H2'	1:CA:383:A:C8	2.48	0.49
6:CG:8:GLN:NE2	6:CG:9:ARG:H	2.10	0.49
9:CJ:46:LYS:HA	9:CJ:68:ARG:HA	1.95	0.49
9:CJ:67:ILE:HG23	9:CJ:67:ILE:O	2.13	0.49
10:CK:69:CYS:O	10:CK:73:VAL:HG13	2.12	0.49
1:CA:1308:U:P	12:CM:97:ARG:HD3	2.53	0.49
19:CT:66:ILE:HG13	19:CT:70:LYS:HG3	1.95	0.49
10:CK:110:THR:CG2	21:CU:4:LYS:HA	2.43	0.49
23:DB:1083:U:C2	23:DB:1086:A:N1	2.81	0.49
23:DB:1142:A:C4	23:DB:1144:A:C8	3.01	0.49
23:DB:1463:C:H2'	23:DB:1464:G:C8	2.48	0.49
23:DB:170:U:H2'	23:DB:171:U:H6	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:208:C:H2'	23:DB:209:C:C6	2.47	0.49
23:DB:242:G:O2'	23:DB:243:U:P	2.70	0.49
23:DB:2484:G:N2	34:DM:118:LYS:HG2	2.27	0.49
23:DB:2809:A:H2'	23:DB:2810:A:H8	1.77	0.49
23:DB:2869:G:H2'	23:DB:2870:C:H6	1.75	0.49
23:DB:312:G:H2'	23:DB:313:G:H8	1.78	0.49
23:DB:318:C:O2'	23:DB:319:G:H5'	2.12	0.49
23:DB:346:A:C8	23:DB:347:A:H1'	2.48	0.49
23:DB:367:G:H2'	23:DB:368:A:O4'	2.12	0.49
23:DB:452:G:N2	23:DB:457:A:O2'	2.46	0.49
23:DB:638:G:H2'	23:DB:639:U:C6	2.48	0.49
28:DF:100:GLU:C	28:DF:102:LEU:H	2.15	0.49
28:DF:41:GLU:O	28:DF:43:ILE:N	2.46	0.49
23:DB:1112:G:H4'	29:DG:1:SER:O	2.13	0.49
31:DJ:54:ILE:HG22	31:DJ:55:ILE:N	2.27	0.49
36:DO:58:ILE:HG13	36:DO:60:GLU:H	1.77	0.49
39:DR:67:GLY:H	39:DR:98:ILE:CA	2.26	0.49
40:DS:3:THR:O	40:DS:4:ILE:HB	2.13	0.49
44:DX:15:ASN:HA	44:DX:17:GLU:OE2	2.13	0.49
1:AA:22:G:H2'	1:AA:23:C:C6	2.47	0.49
1:AA:471:U:H2'	1:AA:472:U:C6	2.47	0.49
1:AA:627:G:H2'	1:AA:628:G:C8	2.48	0.49
2:AC:190:THR:HG22	2:AC:191:THR:H	1.78	0.49
3:AD:150:LYS:HZ3	3:AD:150:LYS:HB2	1.78	0.49
5:AF:54:LEU:HD13	5:AF:55:HIS:N	2.27	0.49
6:AG:139:ASP:HA	6:AG:142:ARG:NH1	2.28	0.49
7:AH:73:SER:HB2	7:AH:129:ALA:HB3	1.94	0.49
1:AA:1343:G:H1'	8:AI:122:ARG:HH12	1.78	0.49
11:AL:54:VAL:HG12	11:AL:56:LEU:HD12	1.95	0.49
13:AN:74:ARG:HH11	13:AN:74:ARG:HG3	1.77	0.49
23:BB:2372:U:OP1	48:B1:44:GLN:HG3	2.13	0.49
49:B2:34:ARG:HG3	49:B2:34:ARG:HH11	1.77	0.49
22:BA:113:C:H2'	22:BA:114:C:C6	2.48	0.49
23:BB:1120:G:O2'	23:BB:1121:C:H5'	2.13	0.49
23:BB:1131:G:H1'	31:BJ:85:LYS:HZ1	1.77	0.49
23:BB:443:A:H2	23:BB:1245:G:N3	2.11	0.49
23:BB:1383:A:H2	23:BB:1405:U:O2	1.95	0.49
23:BB:2354:C:H4'	43:BW:30:VAL:CG2	2.38	0.49
23:BB:2803:G:H2'	23:BB:2804:U:C6	2.47	0.49
23:BB:336:C:O2'	23:BB:337:C:H5'	2.12	0.49
23:BB:678:C:H2'	23:BB:679:C:H6	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:875:G:H2'	23:BB:876:C:O2	2.12	0.49
23:BB:910:A:H2'	23:BB:911:A:C8	2.47	0.49
23:BB:969:G:N2	23:BB:985:C:OP1	2.41	0.49
25:BC:68:ARG:CD	25:BC:127:ASN:HD21	2.26	0.49
26:BD:51:THR:O	26:BD:52:THR:HB	2.13	0.49
26:BD:92:VAL:HG13	26:BD:94:GLN:H	1.78	0.49
27:BE:108:ILE:H	27:BE:108:ILE:CD1	2.26	0.49
27:BE:127:GLU:O	27:BE:157:LEU:HD22	2.13	0.49
28:BF:172:PHE:CD1	28:BF:173:ASP:N	2.80	0.49
30:BH:29:PHE:CD2	30:BH:30:LEU:HD23	2.48	0.49
52:BI:109:ALA:HB1	52:BI:124:MET:CG	2.42	0.49
34:BM:62:LYS:HB3	34:BM:104:GLU:HB2	1.95	0.49
35:BN:65:LEU:CD1	35:BN:68:ALA:HB3	2.43	0.49
35:BN:52:ILE:HG21	35:BN:94:TYR:CE2	2.48	0.49
36:BO:8:ILE:HG22	36:BO:10:ARG:H	1.77	0.49
37:BP:16:VAL:N	37:BP:17:PRO:HD3	2.28	0.49
38:BQ:102:LYS:HG2	38:BQ:103:VAL:H	1.77	0.49
23:BB:583:G:OP1	38:BQ:6:GLY:HA2	2.13	0.49
46:BZ:61:ASN:O	46:BZ:65:ASN:N	2.46	0.49
1:CA:1031:C:H5'	1:CA:1032:G:C8	2.48	0.49
1:CA:1231:G:O2'	1:CA:1232:U:H5'	2.13	0.49
1:CA:1243:C:H2'	1:CA:1244:G:H8	1.78	0.49
1:CA:1356:G:H2'	1:CA:1357:A:C8	2.48	0.49
1:CA:301:G:H2'	1:CA:302:G:C8	2.48	0.49
1:CA:35:G:H2'	1:CA:36:C:H6	1.78	0.49
1:CA:71:A:O2'	1:CA:72:A:H5''	2.13	0.49
1:CA:84:U:O2'	1:CA:86:G:N2	2.46	0.49
2:CC:2:GLN:H	2:CC:2:GLN:NE2	2.10	0.49
9:CJ:15:HIS:HA	9:CJ:18:ILE:HG22	1.94	0.49
1:CA:975:A:H61	9:CJ:50:THR:HB	1.77	0.49
9:CJ:8:ILE:HG13	9:CJ:8:ILE:O	2.13	0.49
12:CM:63:VAL:HG12	12:CM:68:LEU:HB2	1.95	0.49
15:CP:20:VAL:HG21	15:CP:32:PHE:CD2	2.48	0.49
15:CP:3:THR:CG2	15:CP:66:THR:HB	2.39	0.49
16:CQ:30:HIS:CE1	16:CQ:32:ILE:HG22	2.48	0.49
50:D3:4:LYS:CE	50:D3:61:LEU:H	2.26	0.49
22:DA:109:A:H2'	22:DA:110:C:C6	2.48	0.49
23:DB:1439:A:C8	23:DB:1440:U:C6	3.01	0.49
23:DB:1847:A:H4'	23:DB:1848:A:H8	1.78	0.49
23:DB:2737:G:H2'	23:DB:2738:A:C8	2.48	0.49
23:DB:2893:A:H4'	23:DB:2894:G:C5'	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:485:C:O2'	23:DB:486:C:H5'	2.12	0.49
23:DB:518:G:H2'	23:DB:519:U:C6	2.48	0.49
25:DC:12:ARG:HG3	25:DC:19:VAL:O	2.13	0.49
27:DE:138:LEU:HB2	27:DE:143:LEU:HG	1.94	0.49
27:DE:48:THR:HG23	27:DE:85:PHE:H	1.78	0.49
28:DF:8:LYS:HA	28:DF:12:VAL:HG23	1.94	0.49
29:DG:29:ASN:CB	29:DG:78:VAL:HA	2.42	0.49
30:DH:3:VAL:HG21	30:DH:37:VAL:HG21	1.95	0.49
32:DK:104:THR:HB	32:DK:106:GLU:OE1	2.13	0.49
33:DL:75:ALA:HB3	33:DL:108:ALA:HA	1.94	0.49
22:DA:49:C:H4'	36:DO:68:LYS:NZ	2.27	0.49
42:DU:11:ILE:HG22	42:DU:19:GLY:HA2	1.94	0.49
42:DU:28:LEU:HD11	42:DU:31:GLY:CA	2.43	0.49
1:AA:1172:C:H2'	1:AA:1173:U:C6	2.48	0.48
1:AA:970:C:C5	1:AA:1231:G:H1'	2.48	0.48
1:AA:1126:U:O2'	1:AA:1280:A:H2'	2.13	0.48
1:AA:448:A:O2'	1:AA:449:G:H5'	2.13	0.48
1:AA:600:A:H2'	1:AA:601:G:C8	2.48	0.48
1:AA:920:U:H2'	1:AA:921:U:C6	2.48	0.48
3:AD:107:GLY:HA2	3:AD:112:GLU:OE1	2.13	0.48
4:AE:33:THR:O	4:AE:34:ALA:CB	2.61	0.48
8:AI:90:ASP:O	8:AI:93:LEU:HG	2.13	0.48
23:BB:1544:A:H2'	23:BB:1545:A:C8	2.48	0.48
23:BB:1594:U:H2'	23:BB:1595:C:H6	1.75	0.48
23:BB:1747:U:H2'	23:BB:1748:C:C6	2.47	0.48
23:BB:2834:G:H2'	23:BB:2879:A:N6	2.28	0.48
23:BB:435:C:H2'	23:BB:436:C:H5'	1.94	0.48
23:BB:904:G:H2'	23:BB:905:A:H8	1.78	0.48
23:BB:866:A:H61	23:BB:913:U:H4'	1.78	0.48
23:BB:91:A:H2'	23:BB:92:U:C6	2.48	0.48
23:BB:929:U:H2'	23:BB:930:G:C8	2.47	0.48
25:BC:222:THR:C	25:BC:224:MET:H	2.16	0.48
28:BF:135:ILE:O	28:BF:137:PHE:N	2.46	0.48
28:BF:18:GLU:HA	28:BF:18:GLU:OE2	2.12	0.48
29:BG:84:LYS:HA	29:BG:84:LYS:NZ	2.28	0.48
33:BL:85:VAL:HG22	33:BL:98:ALA:HB2	1.94	0.48
36:BO:62:LEU:O	36:BO:64:TYR:N	2.45	0.48
36:BO:8:ILE:HG22	36:BO:10:ARG:CG	2.38	0.48
23:BB:2335:A:P	36:BO:9:ARG:HG3	2.53	0.48
39:BR:33:VAL:HB	39:BR:35:PHE:CE2	2.48	0.48
23:BB:1601:G:OP1	41:BT:62:VAL:HG21	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:99:SER:OG	42:BU:102:ILE:HD11	2.13	0.48
42:BU:46:LYS:N	42:BU:47:PRO:CD	2.72	0.48
23:BB:100:U:C4'	42:BU:90:LYS:HE2	2.26	0.48
1:CA:1004:A:H2'	1:CA:1005:A:O4'	2.13	0.48
1:CA:1275:A:H2'	1:CA:1276:G:C8	2.47	0.48
1:CA:471:U:H2'	1:CA:472:U:C6	2.47	0.48
1:CA:916:U:H2'	1:CA:917:G:H8	1.78	0.48
20:CB:218:ALA:CA	20:CB:221:ARG:HG2	2.41	0.48
4:CE:17:VAL:O	4:CE:17:VAL:HG13	2.12	0.48
5:CF:100:SER:H	17:CR:23:LYS:NZ	2.11	0.48
5:CF:18:VAL:N	5:CF:19:PRO:HD2	2.27	0.48
6:CG:64:ALA:O	6:CG:67:ASN:HB2	2.13	0.48
8:CI:22:PRO:HA	8:CI:60:LEU:HA	1.94	0.48
12:CM:102:LYS:HG3	12:CM:103:THR:H	1.77	0.48
16:CQ:30:HIS:CG	16:CQ:33:TYR:HB2	2.48	0.48
22:DA:49:C:H4'	36:DO:68:LYS:HZ3	1.78	0.48
23:DB:1051:G:H2'	23:DB:1052:C:C6	2.47	0.48
23:DB:1153:C:O2'	23:DB:1154:G:H5'	2.13	0.48
23:DB:1412:U:H2'	23:DB:1413:A:H8	1.76	0.48
23:DB:1583:A:H4'	23:DB:1585:C:N3	2.29	0.48
23:DB:2390:U:O2'	23:DB:2391:G:H5''	2.12	0.48
23:DB:2812:G:H2'	23:DB:2813:A:O4'	2.13	0.48
23:DB:336:C:H5''	42:DU:3:LYS:HZ3	1.76	0.48
23:DB:353:C:O2	23:DB:353:C:H2'	2.13	0.48
23:DB:372:G:C8	46:DZ:57:VAL:HG21	2.48	0.48
23:DB:39:G:H2'	23:DB:40:U:C6	2.48	0.48
23:DB:409:G:H2'	23:DB:410:G:C8	2.48	0.48
23:DB:543:G:H2'	23:DB:545:U:OP1	2.13	0.48
23:DB:685:A:H1'	23:DB:688:U:O4	2.13	0.48
23:DB:833:A:H2'	23:DB:834:G:H8	1.73	0.48
23:DB:899:A:C2	23:DB:900:A:H1'	2.48	0.48
25:DC:168:GLY:C	25:DC:170:TYR:H	2.16	0.48
25:DC:171:VAL:HG23	25:DC:182:LYS:NZ	2.27	0.48
23:DB:1814:G:H5'	25:DC:51:ARG:HG2	1.95	0.48
26:DD:202:ILE:CG2	26:DD:204:LYS:HG2	2.42	0.48
26:DD:34:VAL:HG12	26:DD:91:THR:HG22	1.94	0.48
26:DD:81:GLU:O	26:DD:82:PHE:HB2	2.13	0.48
27:DE:24:ASN:N	27:DE:110:SER:HB2	2.27	0.48
28:DF:39:VAL:CG2	28:DF:84:ILE:HD13	2.42	0.48
28:DF:83:PRO:HG2	28:DF:84:ILE:H	1.78	0.48
29:DG:67:ALA:O	29:DG:71:LEU:HG	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DK:66:LYS:HE3	32:DK:80:ASP:O	2.12	0.48
33:DL:125:LEU:O	33:DL:127:VAL:HG13	2.13	0.48
34:DM:6:ARG:CD	34:DM:71:LYS:HE3	2.42	0.48
37:DP:50:ARG:CG	37:DP:99:LEU:H	2.25	0.48
38:DQ:82:LEU:HG	38:DQ:88:GLU:HG3	1.93	0.48
40:DS:27:LYS:N	40:DS:70:LYS:O	2.46	0.48
43:DW:69:GLU:HG3	43:DW:70:VAL:N	2.24	0.48
43:DW:67:LYS:HZ2	43:DW:70:VAL:HB	1.77	0.48
43:DW:75:ASN:C	43:DW:77:LYS:H	2.16	0.48
44:DX:18:LEU:HA	44:DX:22:LEU:CD1	2.37	0.48
44:DX:18:LEU:HD12	44:DX:47:ARG:NH2	2.28	0.48
46:DZ:25:ARG:HG3	46:DZ:26:SER:N	2.28	0.48
1:AA:1314:C:H2'	1:AA:1315:U:H6	1.77	0.48
1:AA:140:U:H2'	1:AA:141:G:C8	2.48	0.48
1:AA:1510:C:H2'	1:AA:1511:G:H8	1.78	0.48
1:AA:236:A:H2'	1:AA:237:G:H8	1.78	0.48
1:AA:560:A:H4'	1:AA:561:U:H5''	1.94	0.48
1:AA:766:A:C2	1:AA:1525:G:N3	2.81	0.48
1:AA:797:C:OP1	10:AK:125:LYS:HG2	2.13	0.48
1:AA:806:C:H2'	1:AA:807:A:C8	2.47	0.48
1:AA:893:C:H2'	1:AA:894:G:H8	1.78	0.48
1:AA:901:A:H5'	1:AA:902:G:OP2	2.13	0.48
1:AA:940:C:H2'	1:AA:941:G:C8	2.48	0.48
1:AA:405:U:OP2	3:AD:114:ARG:NH2	2.47	0.48
3:AD:54:LEU:C	3:AD:54:LEU:HD13	2.33	0.48
1:AA:546:A:P	3:AD:68:GLU:HB3	2.53	0.48
4:AE:19:ARG:O	4:AE:20:VAL:HB	2.12	0.48
4:AE:90:GLY:O	4:AE:128:GLY:HA3	2.13	0.48
8:AI:9:GLY:HA2	8:AI:80:HIS:CD2	2.48	0.48
9:AJ:88:MET:CE	9:AJ:88:MET:H	2.25	0.48
10:AK:57:SER:O	10:AK:90:PRO:HG2	2.13	0.48
12:AM:56:ARG:O	12:AM:59:VAL:HG12	2.13	0.48
14:AO:80:LEU:HD23	14:AO:80:LEU:C	2.33	0.48
48:B1:31:GLU:O	48:B1:32:LYS:HB2	2.13	0.48
48:B1:5:ARG:O	48:B1:27:ARG:HB2	2.14	0.48
23:BB:1032:A:OP1	51:B4:8:LYS:HB3	2.13	0.48
23:BB:1040:A:H2	23:BB:1115:G:H22	1.59	0.48
23:BB:1137:G:H2'	23:BB:1138:G:O4'	2.13	0.48
23:BB:1287:A:H3'	23:BB:1288:G:N2	2.28	0.48
23:BB:2734:A:C2'	23:BB:2735:G:H5'	2.43	0.48
23:BB:2860:A:H2'	23:BB:2861:U:O4'	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2868:A:H2'	23:BB:2869:G:H8	1.78	0.48
23:BB:2834:G:H2'	23:BB:2879:A:H61	1.78	0.48
23:BB:523:C:H2'	23:BB:524:G:H8	1.78	0.48
23:BB:587:C:H4'	23:BB:588:U:C6	2.48	0.48
23:BB:82:U:H2'	23:BB:83:A:C8	2.48	0.48
23:BB:956:G:H1'	34:BM:81:ARG:HH22	1.77	0.48
23:BB:966:G:H2'	23:BB:967:U:C6	2.47	0.48
26:BD:114:LYS:HB3	35:BN:3:HIS:HE1	1.77	0.48
27:BE:128:ALA:HA	27:BE:157:LEU:CD2	2.42	0.48
27:BE:35:TYR:C	27:BE:37:ALA:H	2.15	0.48
30:BH:9:VAL:C	30:BH:11:ASN:H	2.17	0.48
33:BL:38:GLN:C	33:BL:40:SER:N	2.65	0.48
37:BP:112:ARG:H	37:BP:112:ARG:NH1	2.11	0.48
37:BP:29:VAL:N	37:BP:44:GLY:HA3	2.29	0.48
39:BR:16:GLU:O	39:BR:18:GLN:N	2.45	0.48
44:BX:14:LEU:HD11	44:BX:57:LEU:HD22	1.95	0.48
46:BZ:15:SER:HB3	46:BZ:23:LYS:CD	2.43	0.48
46:BZ:3:LYS:HG2	46:BZ:8:LYS:HA	1.95	0.48
1:CA:1053:G:O2'	1:CA:1199:U:H5	1.96	0.48
9:CJ:41:PRO:O	9:CJ:42:LEU:HB2	2.12	0.48
14:CO:83:ARG:C	14:CO:84:LEU:HD12	2.33	0.48
14:CO:88:ARG:NH1	14:CO:88:ARG:HG3	2.28	0.48
16:CQ:10:ARG:CZ	16:CQ:11:VAL:N	2.76	0.48
18:CS:3:SER:O	18:CS:4:LEU:HD12	2.12	0.48
23:DB:1001:A:H2'	23:DB:1002:G:O4'	2.12	0.48
23:DB:1116:G:H21	34:DM:136:MET:CE	2.25	0.48
23:DB:150:U:H2'	23:DB:151:C:C6	2.48	0.48
23:DB:1790:C:P	25:DC:219:VAL:HB	2.54	0.48
23:DB:1799:G:N7	25:DC:178:GLY:CA	2.72	0.48
23:DB:2488:G:O2'	23:DB:2489:U:H5'	2.12	0.48
23:DB:2893:A:H4'	23:DB:2894:G:O5'	2.12	0.48
23:DB:46:G:H2'	23:DB:47:C:C6	2.48	0.48
23:DB:524:G:H5'	23:DB:539:G:N2	2.27	0.48
23:DB:907:G:O2'	23:DB:908:C:H5'	2.12	0.48
25:DC:124:LYS:HB2	25:DC:125:PRO:HD2	1.93	0.48
25:DC:94:LEU:HD13	25:DC:100:ARG:HG3	1.94	0.48
26:DD:14:ILE:HG23	26:DD:19:GLY:HA3	1.95	0.48
26:DD:46:ARG:H	26:DD:82:PHE:HA	1.78	0.48
27:DE:132:LYS:H	27:DE:134:LEU:CD1	2.26	0.48
27:DE:154:ASP:CG	27:DE:156:ASN:H	2.17	0.48
27:DE:4:VAL:HA	27:DE:14:VAL:HG13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DK:103:VAL:O	32:DK:104:THR:HG23	2.13	0.48
24:DV:29:ILE:HG22	24:DV:90:ASP:HA	1.94	0.48
24:DV:70:ILE:HD12	24:DV:71:LYS:N	2.23	0.48
44:DX:28:LEU:HD13	44:DX:42:LEU:HD21	1.94	0.48
44:DX:43:LEU:N	44:DX:43:LEU:HD22	2.29	0.48
46:DZ:30:HIS:HB3	46:DZ:32:LEU:HD23	1.95	0.48
1:AA:1021:A:H2'	1:AA:1022:A:O4'	2.13	0.48
1:AA:1082:A:O2'	1:AA:1083:U:H5'	2.13	0.48
1:AA:1225:A:H4'	18:AS:77:ARG:HH22	1.77	0.48
1:AA:1247:U:O2'	1:AA:1248:A:H5'	2.13	0.48
1:AA:319:G:O2'	1:AA:320:A:H5'	2.14	0.48
1:AA:408:A:O5'	3:AD:109:THR:HG21	2.12	0.48
1:AA:79:G:H2'	1:AA:80:A:H8	1.74	0.48
2:AC:122:GLN:HE22	2:AC:136:ALA:HB1	1.79	0.48
3:AD:160:LEU:HD22	3:AD:161:ALA:N	2.29	0.48
6:AG:30:MET:HA	6:AG:38:ALA:HB2	1.94	0.48
48:B1:46:VAL:HG22	48:B1:47:ILE:H	1.76	0.48
50:B3:14:LYS:O	50:B3:21:PHE:HB3	2.13	0.48
23:BB:1080:A:O2'	52:BI:126:ARG:HB2	2.13	0.48
23:BB:1230:A:H2'	23:BB:1231:U:H6	1.78	0.48
23:BB:1439:A:C8	23:BB:1440:U:C6	3.01	0.48
23:BB:1463:C:H2'	23:BB:1464:G:C8	2.48	0.48
23:BB:1726:C:H2'	23:BB:1727:C:C6	2.49	0.48
23:BB:231:A:H3'	23:BB:232:G:C8	2.48	0.48
23:BB:2803:G:H2'	23:BB:2804:U:H6	1.79	0.48
23:BB:2893:A:H4'	23:BB:2894:G:O5'	2.12	0.48
23:BB:582:A:H2'	23:BB:583:G:C8	2.47	0.48
23:BB:850:U:H2'	23:BB:851:C:C6	2.48	0.48
23:BB:773:U:H4'	25:BC:47:ARG:HB3	1.94	0.48
26:BD:119:ALA:HB1	26:BD:123:LYS:HB2	1.96	0.48
27:BE:108:ILE:HD12	27:BE:108:ILE:N	2.28	0.48
28:BF:103:ILE:HG13	28:BF:104:THR:N	2.27	0.48
28:BF:133:GLU:HG3	28:BF:134:GLN:H	1.77	0.48
52:BI:79:LEU:HD11	52:BI:131:THR:OG1	2.12	0.48
32:BK:23:LYS:CG	32:BK:24:VAL:N	2.77	0.48
24:BV:76:ASP:CB	34:BM:136:MET:HG3	2.44	0.48
34:BM:76:LYS:HA	34:BM:86:LYS:HD3	1.94	0.48
35:BN:77:ALA:HA	35:BN:80:PHE:CE2	2.48	0.48
36:BO:10:ARG:HH11	36:BO:10:ARG:HG3	1.78	0.48
37:BP:2:ASN:ND2	37:BP:2:ASN:H	2.06	0.48
37:BP:47:ILE:HG23	37:BP:63:ILE:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BQ:71:ASN:HD21	38:BQ:109:VAL:HG11	1.78	0.48
23:BB:1151:A:H4'	38:BQ:80:ASN:OD1	2.12	0.48
39:BR:1:MET:HA	39:BR:46:GLU:HG3	1.94	0.48
39:BR:18:GLN:CB	39:BR:99:THR:HA	2.42	0.48
40:BS:42:LYS:O	40:BS:46:LEU:HG	2.14	0.48
43:BW:16:GLU:HA	43:BW:18:LYS:NZ	2.28	0.48
46:BZ:35:ASP:O	46:BZ:42:PRO:HB3	2.13	0.48
1:CA:1093:A:P	6:CG:3:ARG:HH22	2.36	0.48
1:CA:182:A:H1'	1:CA:183:C:C5	2.48	0.48
1:CA:236:A:H2'	1:CA:237:G:H8	1.78	0.48
1:CA:319:G:O2'	1:CA:320:A:H5'	2.13	0.48
1:CA:513:C:H2'	1:CA:514:C:H6	1.79	0.48
1:CA:505:G:H5'	1:CA:534:U:H2'	1.94	0.48
1:CA:560:A:H4'	1:CA:561:U:H5''	1.94	0.48
1:CA:634:C:O2'	1:CA:635:A:H5'	2.13	0.48
1:CA:8:A:C6	3:CD:205:LYS:HA	2.47	0.48
4:CE:39:GLY:HA3	4:CE:116:VAL:HB	1.95	0.48
15:CP:22:ALA:HB2	15:CP:32:PHE:HA	1.94	0.48
1:CA:255:G:H5'	16:CQ:17:GLU:O	2.13	0.48
19:CT:56:ILE:O	19:CT:60:GLN:HG2	2.13	0.48
48:D1:7:LYS:HB3	48:D1:24:LYS:HZ1	1.77	0.48
49:D2:34:ARG:NH1	49:D2:35:ARG:NH1	2.61	0.48
23:DB:1236:G:C2'	23:DB:1237:A:OP2	2.61	0.48
23:DB:1287:A:H3'	23:DB:1288:G:N2	2.28	0.48
23:DB:197:A:H4'	23:DB:2069:G:OP2	2.12	0.48
23:DB:2378:A:C2'	23:DB:2379:G:H5'	2.43	0.48
23:DB:2063:C:O2	23:DB:2450:A:N1	2.46	0.48
23:DB:2484:G:H1'	34:DM:119:LEU:CD1	2.42	0.48
23:DB:455:C:N4	23:DB:472:A:H2'	2.28	0.48
25:DC:216:ARG:HB3	25:DC:217:PRO:HD2	1.94	0.48
25:DC:261:ARG:HG2	25:DC:261:ARG:NH1	2.28	0.48
25:DC:38:LYS:HG3	25:DC:39:SER:H	1.79	0.48
27:DE:112:LEU:HD12	27:DE:115:GLN:HE22	1.78	0.48
29:DG:84:LYS:HE2	29:DG:163:TYR:CD1	2.48	0.48
23:DB:2751:G:OP2	29:DG:3:VAL:HB	2.14	0.48
31:DJ:121:LYS:CE	31:DJ:121:LYS:N	2.74	0.48
31:DJ:32:LEU:O	31:DJ:36:LEU:HD13	2.13	0.48
32:DK:33:ALA:CB	32:DK:39:ILE:HD11	2.44	0.48
35:DN:97:ILE:HG23	35:DN:113:ILE:HD11	1.95	0.48
35:DN:32:GLU:HA	35:DN:115:LEU:HD21	1.95	0.48
23:DB:2873:A:N3	35:DN:6:SER:HA	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DR:67:GLY:H	39:DR:98:ILE:H	1.60	0.48
39:DR:69:GLY:HA2	39:DR:96:VAL:CA	2.43	0.48
23:DB:974:G:H4'	39:DR:83:TYR:OH	2.13	0.48
43:DW:42:THR:HB	43:DW:75:ASN:ND2	2.28	0.48
1:AA:926:G:N2	1:AA:1505:G:H2'	2.29	0.48
1:AA:418:C:H2'	1:AA:419:C:H6	1.78	0.48
1:AA:554:A:H2'	1:AA:555:U:C6	2.49	0.48
1:AA:708:C:H2'	1:AA:709:U:C6	2.48	0.48
2:AC:66:THR:HA	2:AC:101:ASN:O	2.14	0.48
5:AF:67:PRO:O	5:AF:70:VAL:HG22	2.13	0.48
5:AF:9:MET:HA	5:AF:58:HIS:O	2.13	0.48
10:AK:88:PRO:HD3	21:AU:28:LEU:CD1	2.41	0.48
21:AU:13:VAL:CG1	21:AU:14:ALA:H	2.17	0.48
49:B2:3:ARG:HB3	49:B2:5:PHE:CE1	2.48	0.48
22:BA:52:A:H3'	22:BA:53:A:H8	1.77	0.48
23:BB:1011:G:O2'	23:BB:1013:C:H5''	2.13	0.48
23:BB:1313:U:O2	23:BB:1313:U:H2'	2.12	0.48
23:BB:1312:U:H4'	23:BB:1313:U:O5'	2.12	0.48
23:BB:1407:G:H2'	23:BB:1408:G:H8	1.78	0.48
23:BB:2316:G:H2'	23:BB:2317:A:H8	1.78	0.48
23:BB:2468:A:H2'	23:BB:2476:A:C6	2.48	0.48
23:BB:2544:G:H2'	23:BB:2545:G:H8	1.78	0.48
23:BB:62:U:H3'	23:BB:63:A:C8	2.49	0.48
23:BB:685:A:H1'	23:BB:688:U:O4	2.13	0.48
23:BB:843:G:H2'	23:BB:844:A:C8	2.48	0.48
23:BB:970:U:H2'	23:BB:971:G:H8	1.78	0.48
25:BC:129:LEU:HD22	25:BC:133:ASN:HB2	1.94	0.48
25:BC:155:ARG:HH21	25:BC:155:ARG:HG2	1.79	0.48
25:BC:171:VAL:HA	25:BC:183:VAL:O	2.13	0.48
25:BC:204:LEU:HD11	25:BC:211:ARG:HH11	1.79	0.48
25:BC:216:ARG:HB3	25:BC:217:PRO:CD	2.43	0.48
25:BC:220:ARG:HG3	25:BC:220:ARG:HH11	1.78	0.48
31:BJ:132:HIS:HB3	31:BJ:136:GLN:HE22	1.78	0.48
33:BL:79:LEU:CD1	33:BL:79:LEU:H	2.26	0.48
37:BP:25:VAL:CG2	37:BP:49:ILE:HD11	2.44	0.48
24:BV:57:TYR:HA	24:BV:74:ALA:HB3	1.95	0.48
1:CA:1250:A:H4'	8:CI:69:GLY:N	2.17	0.48
1:CA:1349:A:H1'	1:CA:1374:A:N6	2.28	0.48
1:CA:1411:C:H2'	1:CA:1412:C:C6	2.49	0.48
1:CA:1499:A:O2'	1:CA:1500:A:H5'	2.14	0.48
1:CA:157:U:O2'	1:CA:158:G:H5'	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:399:G:H2'	1:CA:400:C:H6	1.77	0.48
1:CA:448:A:O2'	1:CA:449:G:H5'	2.12	0.48
1:CA:600:A:H2'	1:CA:601:G:C8	2.48	0.48
1:CA:608:A:H2'	1:CA:609:A:O4'	2.14	0.48
1:CA:692:U:H2'	1:CA:694:A:OP2	2.14	0.48
3:CD:202:LEU:C	3:CD:204:SER:H	2.17	0.48
3:CD:25:ARG:HG3	3:CD:25:ARG:O	2.13	0.48
9:CJ:15:HIS:O	9:CJ:18:ILE:HG22	2.13	0.48
1:CA:1309:G:H1'	12:CM:72:ILE:HG13	1.95	0.48
12:CM:89:ARG:CB	12:CM:96:VAL:HG22	2.44	0.48
15:CP:54:LEU:HA	15:CP:57:ILE:CG2	2.43	0.48
16:CQ:45:VAL:HG11	16:CQ:60:ILE:HG21	1.96	0.48
17:CR:62:ARG:C	17:CR:64:LEU:H	2.15	0.48
19:CT:50:PHE:O	19:CT:53:MET:HG3	2.13	0.48
23:DB:1213:A:N1	23:DB:1237:A:H1'	2.27	0.48
23:DB:1244:A:O2'	23:DB:1245:G:H5'	2.13	0.48
23:DB:1268:A:H2'	23:DB:1269:A:O4'	2.13	0.48
23:DB:1312:U:H4'	23:DB:1313:U:O5'	2.12	0.48
23:DB:1511:G:H2'	23:DB:1512:C:H6	1.77	0.48
23:DB:1818:U:OP1	25:DC:155:ARG:HA	2.12	0.48
23:DB:2226:C:H2'	23:DB:2227:A:O4'	2.13	0.48
23:DB:2544:G:H2'	23:DB:2545:G:H8	1.77	0.48
23:DB:2834:G:H2'	23:DB:2879:A:N6	2.28	0.48
23:DB:455:C:H42	23:DB:472:A:H2'	1.77	0.48
23:DB:562:U:C4	23:DB:2036:C:O4'	2.67	0.48
23:DB:856:G:H2'	23:DB:857:G:C8	2.48	0.48
25:DC:225:ASN:N	25:DC:226:PRO:HD3	2.29	0.48
55:DB:3317:HOH:O	26:DD:152:PRO:HG2	2.12	0.48
26:DD:62:LYS:CB	26:DD:63:PRO:HD3	2.43	0.48
27:DE:149:ILE:CD1	27:DE:188:MET:HB2	2.44	0.48
28:DF:56:LEU:CA	28:DF:59:ILE:HG22	2.42	0.48
52:DI:17:ALA:O	52:DI:18:ASN:CB	2.61	0.48
52:DI:70:THR:O	52:DI:70:THR:HG23	2.12	0.48
33:DL:115:GLU:C	33:DL:116:VAL:HG13	2.34	0.48
33:DL:79:LEU:HD11	33:DL:112:LEU:HD23	1.96	0.48
36:DO:29:HIS:HB3	36:DO:36:TYR:HB2	1.95	0.48
40:DS:27:LYS:N	40:DS:27:LYS:HD2	2.29	0.48
41:DT:65:GLY:HA3	41:DT:76:ARG:HH21	1.78	0.48
44:DX:26:PHE:HD1	44:DX:26:PHE:H	1.60	0.48
44:DX:47:ARG:HA	44:DX:50:VAL:HG23	1.95	0.48
45:DY:2:LYS:H	45:DY:37:ARG:HB3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1038:C:H2'	1:AA:1039:G:H8	1.78	0.48
1:AA:1318:A:H4'	18:AS:9:PHE:CD1	2.49	0.48
1:AA:766:A:H2'	1:AA:767:A:O4'	2.13	0.48
1:AA:90:C:H2'	1:AA:91:U:C6	2.47	0.48
15:AP:28:ARG:HG3	15:AP:29:ASN:ND2	2.29	0.48
21:AU:3:ILE:CA	21:AU:19:LYS:HG2	2.35	0.48
22:BA:63:C:H2'	22:BA:64:G:H8	1.79	0.48
23:BB:1299:G:C4'	23:BB:1301:A:H1'	2.40	0.48
23:BB:130:C:O2'	23:BB:131:A:H5'	2.12	0.48
23:BB:142:A:H2'	23:BB:143:C:H6	1.79	0.48
23:BB:1579:A:H2'	23:BB:1580:A:H8	1.78	0.48
23:BB:1599:U:H2'	23:BB:1600:C:C6	2.48	0.48
23:BB:2053:G:OP1	26:BD:149:ASN:O	2.31	0.48
23:BB:2260:C:O2'	23:BB:2261:C:H5'	2.14	0.48
23:BB:2308:G:H3'	23:BB:2308:G:N3	2.28	0.48
23:BB:2722:G:H2'	23:BB:2723:C:C6	2.49	0.48
23:BB:276:U:O2'	23:BB:277:G:H5'	2.14	0.48
23:BB:2812:G:H2'	23:BB:2813:A:O4'	2.13	0.48
23:BB:363:G:H2'	23:BB:364:C:H6	1.78	0.48
23:BB:39:G:O2'	23:BB:40:U:H5'	2.14	0.48
23:BB:572:A:H5''	23:BB:573:U:OP2	2.12	0.48
23:BB:909:A:H2'	23:BB:912:C:H5	1.79	0.48
25:BC:155:ARG:NH2	25:BC:155:ARG:HG2	2.29	0.48
23:BB:1817:G:OP1	25:BC:62:ARG:NH2	2.46	0.48
26:BD:106:LYS:HZ2	26:BD:208:LYS:HD3	1.78	0.48
23:BB:2575:C:C5'	26:BD:149:ASN:N	2.74	0.48
29:BG:174:LYS:HD2	29:BG:174:LYS:N	2.29	0.48
31:BJ:15:TRP:HB3	31:BJ:53:TYR:CE2	2.48	0.48
31:BJ:7:LYS:NZ	31:BJ:49:ASP:H	2.12	0.48
24:BV:76:ASP:HB2	34:BM:136:MET:HE3	1.94	0.48
34:BM:43:ALA:C	34:BM:45:GLN:N	2.67	0.48
38:BQ:88:GLU:HA	39:BR:53:PHE:CG	2.48	0.48
40:BS:32:ALA:O	40:BS:33:LEU:HD13	2.14	0.48
41:BT:6:ARG:HG3	41:BT:10:VAL:HG13	1.95	0.48
41:BT:30:ILE:CG2	41:BT:85:VAL:HG22	2.43	0.48
24:BV:79:ARG:HB3	24:BV:79:ARG:CZ	2.43	0.48
43:BW:39:GLN:C	43:BW:40:ARG:HG3	2.33	0.48
46:BZ:30:HIS:O	46:BZ:48:GLN:NE2	2.47	0.48
46:BZ:50:ASP:CA	46:BZ:53:THR:HG22	2.43	0.48
1:CA:627:G:H2'	1:CA:628:G:H8	1.78	0.48
1:CA:16:A:N1	1:CA:919:A:C2	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:964:A:H2'	1:CA:965:U:H5''	1.95	0.48
2:CC:119:ILE:HD12	2:CC:132:ALA:HB1	1.95	0.48
16:CQ:25:GLU:HA	16:CQ:40:THR:HA	1.96	0.48
16:CQ:80:LYS:H	16:CQ:80:LYS:CD	2.26	0.48
48:D1:4:ILE:HG23	48:D1:4:ILE:O	2.13	0.48
23:DB:1178:C:C2	23:DB:1179:G:N7	2.81	0.48
23:DB:1299:G:H4'	23:DB:1301:A:C1'	2.42	0.48
23:DB:1322:A:C2'	23:DB:1323:C:H5'	2.44	0.48
23:DB:1581:G:O2'	23:DB:1582:C:H5'	2.13	0.48
23:DB:1726:C:H2'	23:DB:1727:C:C6	2.49	0.48
23:DB:2563:U:H2'	23:DB:2565:A:OP2	2.13	0.48
23:DB:2568:U:H2'	23:DB:2569:G:O4'	2.14	0.48
23:DB:2803:G:H2'	23:DB:2804:U:H6	1.79	0.48
23:DB:2840:C:H2'	23:DB:2841:C:C6	2.48	0.48
23:DB:673:C:O2'	23:DB:674:G:H5'	2.13	0.48
23:DB:723:C:H2'	23:DB:724:U:H6	1.79	0.48
25:DC:140:VAL:HG11	25:DC:163:ILE:CD1	2.42	0.48
27:DE:147:LEU:N	27:DE:147:LEU:HD12	2.27	0.48
27:DE:152:GLU:OE2	27:DE:188:MET:HA	2.13	0.48
27:DE:15:SER:CB	27:DE:196:VAL:HG22	2.41	0.48
30:DH:90:LEU:HD13	30:DH:122:LEU:O	2.14	0.48
30:DH:135:HIS:CG	30:DH:136:SER:N	2.81	0.48
31:DJ:114:LEU:HG	31:DJ:118:MET:CE	2.43	0.48
37:DP:48:ALA:HB3	37:DP:64:SER:HB3	1.94	0.48
41:DT:2:ILE:HG12	41:DT:3:ARG:N	2.27	0.48
42:DU:96:LYS:O	42:DU:97:SER:HB3	2.13	0.48
43:DW:44:PHE:HB3	43:DW:77:LYS:C	2.34	0.48
46:DZ:62:LYS:CA	46:DZ:65:ASN:HD21	2.26	0.48
1:AA:999:C:O2'	1:AA:1000:A:H5'	2.13	0.48
1:AA:1025:U:H5''	1:AA:1026:G:OP1	2.14	0.48
1:AA:1149:C:H2'	1:AA:1150:A:C8	2.48	0.48
1:AA:978:A:O2'	1:AA:1322:C:H5	1.96	0.48
1:AA:429:U:H3'	3:AD:8:LEU:HD23	1.96	0.48
1:AA:869:G:H4'	1:AA:872:A:C8	2.49	0.48
2:AC:90:VAL:HA	2:AC:93:ILE:HG21	1.95	0.48
14:AO:54:GLY:O	14:AO:58:MET:HG2	2.13	0.48
18:AS:46:LEU:O	18:AS:60:PHE:HA	2.12	0.48
18:AS:4:LEU:CA	18:AS:5:LYS:HE3	2.43	0.48
47:B0:36:LYS:HZ3	47:B0:41:HIS:HA	1.78	0.48
23:BB:1017:G:H2'	23:BB:1018:U:C6	2.49	0.48
23:BB:1079:C:O2'	52:BI:133:ARG:NH2	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1332:G:H2'	23:BB:1332:G:N3	2.29	0.48
23:BB:1355:G:O2'	23:BB:1356:G:H5'	2.14	0.48
23:BB:1688:U:O2	23:BB:1700:A:H5'	2.13	0.48
23:BB:2092:U:H4'	23:BB:2093:G:O5'	2.14	0.48
23:BB:2180:U:H2'	23:BB:2181:U:C5	2.48	0.48
23:BB:2307:G:N3	23:BB:2307:G:H2'	2.29	0.48
23:BB:231:A:H3'	23:BB:232:G:H8	1.79	0.48
23:BB:2530:A:H61	29:BG:155:PRO:HG3	1.78	0.48
23:BB:2593:U:O2'	23:BB:2594:C:H5'	2.13	0.48
23:BB:2645:G:H4'	23:BB:2732:G:H2'	1.96	0.48
23:BB:2776:A:H4'	23:BB:2777:G:O5'	2.13	0.48
23:BB:306:U:H2'	23:BB:307:G:O4'	2.13	0.48
23:BB:720:U:H2'	23:BB:721:A:H8	1.78	0.48
25:BC:71:ASP:O	25:BC:73:ILE:N	2.46	0.48
26:BD:142:VAL:C	26:BD:144:GLY:N	2.67	0.48
26:BD:2:ILE:HG22	26:BD:204:LYS:HG3	1.95	0.48
29:BG:114:HIS:HB2	29:BG:150:TYR:HE2	1.79	0.48
30:BH:7:ASP:O	30:BH:9:VAL:HG23	2.13	0.48
30:BH:7:ASP:OD1	30:BH:35:LYS:HB3	2.14	0.48
34:BM:73:ILE:CD1	34:BM:92:TRP:H	2.25	0.48
35:BN:16:HIS:O	35:BN:20:MET:HG2	2.13	0.48
35:BN:2:ARG:O	35:BN:3:HIS:CB	2.62	0.48
35:BN:7:GLY:O	35:BN:8:ARG:CB	2.61	0.48
37:BP:50:ARG:NH1	37:BP:62:LYS:HB3	2.29	0.48
37:BP:52:ARG:N	37:BP:60:VAL:HG21	2.29	0.48
39:BR:45:GLU:HA	39:BR:50:GLY:HA2	1.95	0.48
39:BR:41:ILE:O	39:BR:54:VAL:HG22	2.14	0.48
39:BR:71:LYS:O	39:BR:94:THR:HA	2.13	0.48
24:BV:9:ARG:HB2	24:BV:39:ALA:O	2.13	0.48
45:BY:18:LYS:HD2	45:BY:53:MET:CE	2.42	0.48
1:CA:123:U:H5''	1:CA:311:C:O2'	2.12	0.48
1:CA:140:U:H2'	1:CA:141:G:C8	2.48	0.48
1:CA:218:U:H2'	1:CA:219:U:H6	1.78	0.48
1:CA:280:C:O2	16:CQ:39:ARG:HG3	2.13	0.48
1:CA:840:C:C2	1:CA:842:U:H4'	2.49	0.48
20:CB:95:TRP:HZ3	20:CB:98:GLY:H	1.62	0.48
1:CA:1240:U:P	6:CG:115:MET:H	2.37	0.48
6:CG:55:LYS:HG3	6:CG:56:SER:N	2.28	0.48
6:CG:59:GLU:HA	6:CG:62:GLU:CD	2.34	0.48
8:CI:112:ARG:HD2	13:CN:100:TRP:OXT	2.13	0.48
10:CK:30:ILE:HG22	10:CK:45:THR:HB	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:CO:10:ILE:HD11	14:CO:29:ALA:HB1	1.94	0.48
14:CO:84:LEU:HB3	14:CO:86:LEU:HD13	1.96	0.48
21:CU:15:LEU:HA	21:CU:17:ARG:NH1	2.24	0.48
51:D4:30:GLU:CB	51:D4:33:HIS:HB2	2.41	0.48
51:D4:24:ARG:HE	51:D4:37:GLN:CA	2.26	0.48
22:DA:94:A:OP1	24:DV:19:ARG:HD3	2.13	0.48
23:DB:1313:U:O2	23:DB:1313:U:H2'	2.12	0.48
23:DB:1791:A:H61	23:DB:1828:G:H1'	1.78	0.48
23:DB:2000:C:O2'	23:DB:2001:C:H5'	2.14	0.48
23:DB:2026:U:H2'	23:DB:2027:G:H8	1.77	0.48
23:DB:2512:C:OP1	26:DD:128:ARG:N	2.47	0.48
23:DB:2516:A:O2'	23:DB:2517:C:H5'	2.13	0.48
23:DB:274:C:H2'	23:DB:275:C:O4'	2.12	0.48
23:DB:2862:G:H2'	23:DB:2863:C:H6	1.79	0.48
23:DB:654:A:H2'	23:DB:655:A:C5'	2.33	0.48
23:DB:859:G:O2'	23:DB:860:U:OP2	2.32	0.48
25:DC:77:VAL:HG21	25:DC:109:LEU:HD23	1.95	0.48
25:DC:53:ILE:HG23	25:DC:216:ARG:HB2	1.95	0.48
26:DD:153:GLY:C	26:DD:155:VAL:N	2.65	0.48
26:DD:175:LEU:HD21	26:DD:192:ALA:HB3	1.95	0.48
23:DB:2633:G:C1'	26:DD:62:LYS:HG3	2.39	0.48
27:DE:56:GLY:O	27:DE:57:LYS:HG3	2.14	0.48
28:DF:136:ILE:O	28:DF:138:PRO:HD3	2.13	0.48
33:DL:37:GLY:O	33:DL:38:GLN:HG3	2.14	0.48
33:DL:50:PHE:O	33:DL:51:GLU:C	2.51	0.48
34:DM:83:GLY:O	34:DM:84:LYS:HB2	2.12	0.48
34:DM:43:ALA:HB3	34:DM:91:TYR:CE2	2.48	0.48
40:DS:72:THR:HG22	40:DS:106:VAL:O	2.13	0.48
42:DU:48:VAL:HG13	42:DU:51:LEU:N	2.29	0.48
24:DV:77:VAL:HG13	24:DV:89:ILE:CD1	2.44	0.48
43:DW:20:LEU:CD1	43:DW:31:LEU:HB2	2.44	0.48
43:DW:43:LYS:O	43:DW:44:PHE:CB	2.60	0.48
43:DW:67:LYS:CG	43:DW:71:LYS:HB2	2.36	0.48
43:DW:73:PRO:C	43:DW:74:LYS:HD2	2.34	0.48
44:DX:55:THR:C	44:DX:57:LEU:H	2.15	0.48
1:AA:1352:C:H2'	1:AA:1353:G:C8	2.49	0.48
1:AA:147:G:H2'	1:AA:148:G:C8	2.48	0.48
1:AA:340:U:H2'	1:AA:341:C:H6	1.79	0.48
1:AA:397:A:H5'	1:AA:398:U:OP1	2.13	0.48
1:AA:736:C:H2'	1:AA:737:C:H6	1.77	0.48
1:AA:79:G:H2'	1:AA:80:A:O4'	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:922:G:H2'	1:AA:923:A:H8	1.78	0.48
1:AA:948:C:O2'	1:AA:949:A:H5'	2.14	0.48
20:AB:68:PHE:HA	20:AB:161:PHE:O	2.12	0.48
3:AD:2:ARG:HG3	3:AD:114:ARG:NH1	2.29	0.48
10:AK:80:ASN:HD21	10:AK:107:THR:CG2	2.27	0.48
11:AL:113:ARG:HA	11:AL:118:VAL:HG23	1.96	0.48
19:AT:19:HIS:O	19:AT:23:ARG:HG2	2.13	0.48
47:B0:11:LYS:C	47:B0:13:GLY:H	2.15	0.48
33:BL:60:ARG:NH1	50:B3:11:LYS:HE2	2.27	0.48
23:BB:1561:C:H2'	23:BB:1562:U:C6	2.49	0.48
23:BB:1759:A:H4'	23:BB:2715:C:O4'	2.13	0.48
23:BB:2000:C:O2'	23:BB:2001:C:H5'	2.14	0.48
23:BB:2743:U:C3'	23:BB:2744:G:H5''	2.44	0.48
23:BB:2795:C:O5'	23:BB:2795:C:H6	1.97	0.48
23:BB:67:U:C4	23:BB:68:G:N7	2.82	0.48
23:BB:712:G:H2'	23:BB:713:G:O4'	2.13	0.48
25:BC:171:VAL:HG21	25:BC:182:LYS:HD3	1.96	0.48
25:BC:143:VAL:HG11	25:BC:173:LEU:HD11	1.95	0.48
25:BC:218:THR:HG22	25:BC:218:THR:O	2.13	0.48
25:BC:226:PRO:HD3	25:BC:233:GLY:N	2.29	0.48
25:BC:51:ARG:HD3	25:BC:51:ARG:O	2.13	0.48
26:BD:34:VAL:HG11	26:BD:71:ALA:CB	2.44	0.48
26:BD:46:ARG:HD2	26:BD:80:TRP:CE2	2.49	0.48
30:BH:19:VAL:HG12	30:BH:20:ASN:N	2.28	0.48
52:BI:75:ALA:HB2	52:BI:112:LYS:HE2	1.95	0.48
23:BB:1059:G:N2	52:BI:130:GLY:HA3	2.29	0.48
31:BJ:25:LEU:HD12	31:BJ:62:VAL:HG12	1.94	0.48
31:BJ:7:LYS:HZ3	31:BJ:48:VAL:CA	2.27	0.48
33:BL:110:VAL:HG23	33:BL:126:ARG:CB	2.43	0.48
34:BM:20:LEU:CB	34:BM:38:ARG:HD3	2.43	0.48
35:BN:102:PHE:N	35:BN:102:PHE:HD2	2.11	0.48
36:BO:58:ILE:HG13	36:BO:63:LYS:HB3	1.94	0.48
40:BS:28:LYS:HB3	40:BS:71:VAL:HG21	1.96	0.48
45:BY:9:THR:HG23	45:BY:10:ARG:N	2.25	0.48
1:CA:1116:U:O2'	1:CA:1117:A:H5'	2.13	0.48
1:CA:1164:G:O2'	1:CA:1165:U:H5'	2.13	0.48
1:CA:1287:A:H2'	1:CA:1288:A:H8	1.76	0.48
1:CA:1390:U:H2'	1:CA:1391:U:C6	2.49	0.48
1:CA:1516:G:H2'	1:CA:1518:A:OP2	2.13	0.48
1:CA:659:U:H2'	1:CA:660:C:C6	2.48	0.48
1:CA:893:C:H2'	1:CA:894:G:H8	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	2.48	0.48
20:CB:153:MET:C	20:CB:155:GLY:H	2.17	0.48
3:CD:27:ILE:C	3:CD:29:THR:H	2.16	0.48
4:CE:148:SER:HB2	4:CE:149:PRO:HD2	1.94	0.48
6:CG:139:ASP:O	6:CG:142:ARG:HB3	2.14	0.48
8:CI:33:SER:HB2	8:CI:36:GLN:CG	2.44	0.48
9:CJ:37:ARG:HG2	9:CJ:37:ARG:NH1	2.29	0.48
9:CJ:64:GLN:CB	13:CN:98:ALA:HB3	2.43	0.48
10:CK:70:ALA:N	10:CK:73:VAL:HG22	2.29	0.48
13:CN:40:ARG:CZ	18:CS:6:LYS:HB2	2.44	0.48
16:CQ:60:ILE:HG12	16:CQ:72:TRP:HE3	1.78	0.48
18:CS:56:HIS:CD2	18:CS:56:HIS:N	2.80	0.48
23:DB:1407:G:H2'	23:DB:1408:G:H8	1.78	0.48
23:DB:1727:C:H2'	23:DB:1728:C:H6	1.74	0.48
23:DB:2105:U:H2'	23:DB:2106:U:C6	2.48	0.48
23:DB:2800:A:H2'	23:DB:2801:G:C4'	2.43	0.48
23:DB:417:C:H2'	23:DB:418:C:H6	1.78	0.48
23:DB:493:G:O2'	23:DB:494:G:H5'	2.14	0.48
23:DB:851:C:H2'	23:DB:852:U:C6	2.49	0.48
23:DB:955:U:H5'	23:DB:956:G:OP2	2.12	0.48
25:DC:153:LEU:C	25:DC:155:ARG:H	2.17	0.48
25:DC:212:TRP:CZ3	25:DC:217:PRO:HD3	2.49	0.48
26:DD:48:ILE:CG2	26:DD:49:GLN:N	2.76	0.48
30:DH:2:GLN:CB	30:DH:19:VAL:HA	2.43	0.48
52:DI:90:GLY:C	52:DI:92:PRO:HD3	2.34	0.48
31:DJ:68:LYS:HD2	31:DJ:72:LYS:CB	2.43	0.48
35:DN:2:ARG:NH1	35:DN:4:ARG:HH12	2.11	0.48
36:DO:4:LYS:O	36:DO:7:ARG:HG2	2.12	0.48
37:DP:85:VAL:O	37:DP:87:ARG:NE	2.46	0.48
38:DQ:50:ARG:HH22	38:DQ:53:LYS:NZ	2.12	0.48
42:DU:22:GLY:HA3	42:DU:35:VAL:CG2	2.43	0.48
44:DX:16:THR:HG1	44:DX:19:LEU:HB3	1.79	0.48
45:DY:3:THR:CA	45:DY:37:ARG:H	2.23	0.48
1:AA:1037:C:H2'	1:AA:1038:C:C6	2.48	0.48
1:AA:1225:A:N3	1:AA:1225:A:H2'	2.29	0.48
1:AA:1487:G:O2'	1:AA:1488:G:H5'	2.13	0.48
1:AA:267:C:OP2	16:AQ:68:LYS:HD2	2.14	0.48
1:AA:363:A:OP2	11:AL:30:ARG:NH1	2.47	0.48
1:AA:591:U:H2'	1:AA:592:G:C8	2.48	0.48
3:AD:66:VAL:HG13	3:AD:70:GLN:OE1	2.13	0.48
3:AD:201:GLU:OE1	4:AE:104:ILE:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AN:50:LEU:CD2	13:AN:51:PRO:HD3	2.41	0.48
16:AQ:17:GLU:C	16:AQ:19:SER:H	2.17	0.48
16:AQ:30:HIS:H	16:AQ:35:LYS:H	1.61	0.48
16:AQ:28:VAL:C	16:AQ:36:PHE:HA	2.33	0.48
50:B3:12:ARG:NH2	50:B3:55:GLY:HA3	2.28	0.48
23:BB:1206:G:H2'	23:BB:1207:C:C6	2.49	0.48
23:BB:132:G:H2'	23:BB:133:U:C6	2.49	0.48
23:BB:1467:U:C4	23:BB:1546:G:C2	3.01	0.48
23:BB:1583:A:H4'	23:BB:1585:C:N3	2.28	0.48
23:BB:1657:U:P	26:BD:141:ARG:HB2	2.54	0.48
23:BB:1791:A:H61	23:BB:1828:G:H1'	1.78	0.48
23:BB:1841:U:H2'	23:BB:1842:G:H8	1.79	0.48
23:BB:188:G:O2'	23:BB:189:G:H5'	2.14	0.48
23:BB:207:A:H2'	23:BB:208:C:O4'	2.14	0.48
23:BB:2149:U:O2'	23:BB:2150:C:H5'	2.13	0.48
23:BB:2259:U:O2	23:BB:2259:U:H2'	2.13	0.48
23:BB:2394:C:OP1	33:BL:63:LYS:O	2.30	0.48
23:BB:2893:A:H4'	23:BB:2894:G:C5'	2.43	0.48
23:BB:673:C:O2'	23:BB:674:G:H5'	2.13	0.48
23:BB:942:G:O2'	23:BB:943:A:H5'	2.14	0.48
23:BB:1818:U:H5''	25:BC:155:ARG:HG2	1.95	0.48
25:BC:229:HIS:ND1	25:BC:230:PRO:HD2	2.28	0.48
25:BC:87:SER:N	25:BC:155:ARG:HH12	2.11	0.48
23:BB:2572:A:OP2	26:BD:151:THR:HA	2.14	0.48
26:BD:1:MET:SD	26:BD:84:LEU:HD13	2.54	0.48
28:BF:59:ILE:HD11	28:BF:137:PHE:CG	2.49	0.48
29:BG:18:ILE:N	29:BG:18:ILE:HD13	2.29	0.48
30:BH:81:ALA:CB	30:BH:147:VAL:H	2.27	0.48
52:BI:58:ILE:N	52:BI:58:ILE:HD12	2.29	0.48
52:BI:63:ASP:C	52:BI:65:SER:N	2.67	0.48
31:BJ:21:THR:HA	31:BJ:61:LYS:HB2	1.95	0.48
31:BJ:7:LYS:HZ3	31:BJ:49:ASP:H	1.61	0.48
33:BL:38:GLN:H	33:BL:41:ARG:NH1	2.11	0.48
35:BN:99:LYS:HZ3	35:BN:99:LYS:HB3	1.78	0.48
37:BP:93:LYS:CB	37:BP:96:LEU:HA	2.43	0.48
23:BB:1156:A:H62	38:BQ:47:ARG:CD	2.27	0.48
39:BR:37:GLU:HG2	39:BR:62:GLU:N	2.19	0.48
41:BT:34:VAL:HG11	41:BT:43:ILE:HD11	1.94	0.48
41:BT:6:ARG:C	41:BT:8:LEU:N	2.67	0.48
24:BV:26:PHE:HZ	24:BV:47:VAL:HB	1.79	0.48
24:BV:77:VAL:HG13	24:BV:77:VAL:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BV:78:GLN:HB2	24:BV:88:HIS:O	2.14	0.48
46:BZ:31:ASP:O	46:BZ:32:LEU:HB3	2.14	0.48
1:CA:1008:U:H3'	1:CA:1009:U:H5''	1.95	0.48
1:CA:1020:G:N3	1:CA:1020:G:H2'	2.29	0.48
1:CA:1260:G:HO2'	1:CA:1261:A:H8	1.59	0.48
1:CA:1310:G:O2'	1:CA:1311:A:H5'	2.13	0.48
1:CA:1435:G:H2'	1:CA:1436:U:C6	2.49	0.48
1:CA:591:U:H2'	1:CA:592:G:C8	2.48	0.48
1:CA:766:A:H2'	1:CA:767:A:O4'	2.13	0.48
20:CB:86:CYS:SG	20:CB:87:ASP:N	2.86	0.48
3:CD:156:ALA:O	3:CD:160:LEU:HD12	2.14	0.48
4:CE:17:VAL:HA	4:CE:33:THR:O	2.13	0.48
4:CE:19:ARG:O	4:CE:20:VAL:HB	2.14	0.48
6:CG:112:ASP:HB2	6:CG:118:ARG:CG	2.41	0.48
8:CI:40:ARG:HA	8:CI:44:ARG:NH1	2.29	0.48
8:CI:56:MET:C	8:CI:58:GLU:H	2.17	0.48
10:CK:121:ARG:HH21	21:CU:34:ARG:NE	2.11	0.48
11:CL:75:GLU:OE2	11:CL:76:HIS:N	2.47	0.48
12:CM:18:LEU:HD23	12:CM:24:VAL:HG21	1.94	0.48
17:CR:20:ILE:HD12	17:CR:53:GLN:HE22	1.79	0.48
21:CU:6:ARG:C	21:CU:7:GLU:HG2	2.34	0.48
23:DB:1666:G:O2'	23:DB:1667:G:H5'	2.13	0.48
23:DB:1706:C:C2	23:DB:1757:A:H5'	2.48	0.48
23:DB:1779:U:OP2	23:DB:1784:A:N6	2.46	0.48
23:DB:2106:U:C2'	23:DB:2107:G:OP1	2.62	0.48
23:DB:2344:U:H4'	23:DB:2345:G:OP1	2.13	0.48
23:DB:197:A:H62	23:DB:2430:A:H2'	1.78	0.48
23:DB:2526:G:H2'	23:DB:2527:C:C6	2.48	0.48
23:DB:810:U:N3	33:DL:37:GLY:HA2	2.29	0.48
23:DB:871:U:H4'	34:DM:68:PHE:CE1	2.49	0.48
23:DB:951:C:O2'	23:DB:952:G:H5'	2.14	0.48
25:DC:139:THR:O	25:DC:191:LEU:HD12	2.14	0.48
25:DC:267:VAL:HG12	25:DC:268:ARG:N	2.28	0.48
23:DB:1098:A:N3	52:DI:3:LYS:O	2.46	0.48
31:DJ:105:VAL:HA	31:DJ:108:MET:HE3	1.94	0.48
31:DJ:114:LEU:HG	31:DJ:118:MET:HE1	1.96	0.48
35:DN:14:SER:HA	35:DN:17:ARG:HH12	1.79	0.48
37:DP:18:SER:HB2	37:DP:87:ARG:NH2	2.28	0.48
38:DQ:98:ALA:HA	38:DQ:105:PHE:CG	2.49	0.48
39:DR:32:THR:HB	39:DR:65:ALA:C	2.34	0.48
41:DT:47:VAL:HG13	41:DT:48:GLN:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DW:67:LYS:HE3	43:DW:71:LYS:H	1.78	0.48
1:AA:1018:G:H2'	1:AA:1019:A:H8	1.77	0.48
1:AA:1071:C:O2'	1:AA:1072:G:H5'	2.14	0.48
1:AA:476:U:O2'	1:AA:477:C:H5'	2.14	0.48
1:AA:608:A:H2'	1:AA:609:A:O4'	2.14	0.48
1:AA:66:A:H8	1:AA:66:A:O5'	1.97	0.48
1:AA:840:C:C2	1:AA:842:U:H4'	2.49	0.48
2:AC:154:GLY:HA2	2:AC:163:ARG:O	2.12	0.48
2:AC:57:GLU:HB2	2:AC:64:ARG:CB	2.43	0.48
3:AD:37:PRO:HD2	3:AD:41:GLY:HA3	1.96	0.48
3:AD:58:GLN:O	3:AD:62:ARG:HG2	2.12	0.48
4:AE:89:THR:HG22	4:AE:90:GLY:N	2.29	0.48
8:AI:51:LEU:HD11	8:AI:82:ILE:HG21	1.96	0.48
8:AI:79:ARG:O	8:AI:83:THR:HG22	2.14	0.48
11:AL:78:VAL:O	11:AL:102:ASP:HB2	2.14	0.48
13:AN:33:VAL:HG23	13:AN:40:ARG:HH21	1.79	0.48
18:AS:66:VAL:C	18:AS:68:HIS:H	2.17	0.48
47:B0:28:SER:O	47:B0:36:LYS:HB2	2.13	0.48
23:BB:1157:G:H2'	23:BB:1158:C:H6	1.79	0.48
23:BB:839:U:O2'	23:BB:1191:G:H1'	2.14	0.48
23:BB:1433:A:H2'	23:BB:1434:A:O4'	2.13	0.48
23:BB:2059:A:N6	23:BB:2503:A:H2'	2.29	0.48
23:BB:2291:U:OP1	23:BB:2380:C:O2'	2.31	0.48
23:BB:2485:G:H1'	34:BM:118:LYS:HZ3	1.79	0.48
23:BB:2683:C:H2'	23:BB:2684:U:C6	2.49	0.48
23:BB:545:U:H2'	23:BB:547:A:P	2.54	0.48
23:BB:562:U:H2'	23:BB:572:A:O4'	2.14	0.48
23:BB:633:A:H2'	23:BB:634:C:C5'	2.41	0.48
23:BB:728:G:H4'	25:BC:12:ARG:HH22	1.79	0.48
25:BC:164:VAL:HB	25:BC:167:ASP:OD2	2.13	0.48
25:BC:231:HIS:ND1	25:BC:231:HIS:O	2.47	0.48
28:BF:55:ASP:CG	28:BF:148:VAL:HG11	2.34	0.48
29:BG:70:LEU:O	29:BG:74:MET:HG3	2.14	0.48
30:BH:3:VAL:HG22	30:BH:21:VAL:HG21	1.96	0.48
52:BI:17:ALA:O	52:BI:18:ASN:CB	2.62	0.48
31:BJ:35:ARG:NE	31:BJ:39:LYS:HB3	2.24	0.48
31:BJ:46:PRO:O	31:BJ:47:HIS:HB2	2.14	0.48
35:BN:9:GLN:HE21	35:BN:17:ARG:HH12	1.61	0.48
37:BP:29:VAL:C	37:BP:44:GLY:HA3	2.34	0.48
38:BQ:35:PHE:HB3	38:BQ:39:ILE:HD12	1.96	0.48
38:BQ:52:ARG:CZ	38:BQ:55:GLN:HE21	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:BR:66:HIS:N	39:BR:98:ILE:HD13	2.28	0.48
40:BS:29:VAL:C	40:BS:31:GLN:H	2.16	0.48
41:BT:20:ALA:C	41:BT:22:THR:H	2.17	0.48
41:BT:74:ILE:HG22	41:BT:75:GLY:N	2.29	0.48
42:BU:96:LYS:HE2	42:BU:98:ASN:HB3	1.96	0.48
42:BU:95:PHE:HD1	42:BU:96:LYS:HG2	1.78	0.48
45:BY:35:VAL:CG1	45:BY:37:ARG:HH21	2.27	0.48
1:CA:1254:A:H2'	1:CA:1255:G:C8	2.49	0.48
1:CA:1366:C:H2'	1:CA:1367:C:H6	1.77	0.48
1:CA:554:A:H2'	1:CA:555:U:C6	2.49	0.48
1:CA:869:G:H4'	1:CA:872:A:C8	2.49	0.48
2:CC:168:ARG:HG2	2:CC:169:GLU:N	2.28	0.48
5:CF:12:PRO:HB3	5:CF:56:LYS:O	2.13	0.48
5:CF:38:ARG:NH2	5:CF:63:ASN:ND2	2.59	0.48
6:CG:75:LYS:HE3	6:CG:76:SER:O	2.13	0.48
7:CH:88:LYS:HD3	7:CH:89:ASP:N	2.29	0.48
8:CI:29:ILE:HB	8:CI:64:ILE:HD12	1.95	0.48
10:CK:110:THR:HG22	21:CU:5:VAL:N	2.23	0.48
10:CK:48:GLY:C	10:CK:50:GLY:H	2.17	0.48
11:CL:51:VAL:HG12	11:CL:52:CYS:N	2.27	0.48
12:CM:102:LYS:NZ	12:CM:103:THR:HG23	2.28	0.48
15:CP:54:LEU:HD13	15:CP:80:LYS:HZ2	1.78	0.48
16:CQ:10:ARG:NE	16:CQ:56:ASP:O	2.44	0.48
17:CR:47:ARG:CZ	17:CR:48:ALA:H	2.27	0.48
19:CT:2:ASN:O	19:CT:3:ILE:C	2.52	0.48
21:CU:39:LYS:H	21:CU:40:PRO:CD	2.21	0.48
50:D3:4:LYS:HE3	50:D3:61:LEU:CB	2.44	0.48
23:DB:1197:G:O2'	23:DB:1198:U:H5'	2.14	0.48
23:DB:1213:A:C6	23:DB:1237:A:H1'	2.49	0.48
23:DB:1438:U:O2'	23:DB:1439:A:H5'	2.14	0.48
23:DB:1467:U:C4	23:DB:1546:G:C2	3.01	0.48
23:DB:150:U:H2'	23:DB:151:C:H6	1.78	0.48
23:DB:1688:U:O2	23:DB:1700:A:H5'	2.13	0.48
23:DB:1819:A:OP1	25:DC:155:ARG:HB3	2.14	0.48
23:DB:1821:A:H2'	23:DB:1822:C:H6	1.76	0.48
23:DB:2038:G:H2'	23:DB:2039:U:C6	2.49	0.48
23:DB:2466:C:O2'	34:DM:118:LYS:HE2	2.13	0.48
23:DB:2773:C:H5''	26:DD:169:ARG:HB2	1.95	0.48
23:DB:704:G:HO2'	23:DB:705:A:P	2.37	0.48
23:DB:963:U:H2'	23:DB:964:C:C6	2.49	0.48
25:DC:128:THR:HA	25:DC:189:ALA:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:137:GLY:C	25:DC:139:THR:N	2.66	0.48
26:DD:121:THR:HG21	26:DD:143:PRO:HD3	1.96	0.48
27:DE:134:LEU:HD22	27:DE:134:LEU:N	2.27	0.48
27:DE:14:VAL:HG11	27:DE:16:GLU:CD	2.34	0.48
33:DL:77:ILE:HG21	33:DL:92:LEU:HG	1.96	0.48
36:DO:39:VAL:HG13	36:DO:39:VAL:O	2.13	0.48
37:DP:87:ARG:HH11	37:DP:87:ARG:HG2	1.79	0.48
38:DQ:55:GLN:O	38:DQ:58:GLN:HG3	2.12	0.48
41:DT:63:VAL:O	41:DT:79:ASP:HA	2.14	0.48
23:DB:2261:C:OP2	43:DW:13:ARG:HB2	2.14	0.48
45:DY:2:LYS:CA	45:DY:37:ARG:HB2	2.44	0.48
1:AA:110:C:H2'	1:AA:111:G:O4'	2.14	0.48
1:AA:113:G:H1'	1:AA:354:G:H5'	1.96	0.48
1:AA:333:U:H2'	1:AA:334:C:C6	2.49	0.48
1:AA:36:C:H5''	11:AL:119:LYS:HB3	1.94	0.48
1:AA:659:U:H2'	1:AA:660:C:C6	2.49	0.48
1:AA:16:A:N1	1:AA:919:A:H2	2.12	0.48
20:AB:127:LYS:HB2	20:AB:127:LYS:NZ	2.29	0.48
3:AD:154:VAL:HG13	3:AD:155:LYS:N	2.25	0.48
12:AM:18:LEU:O	12:AM:21:ILE:HG13	2.13	0.48
19:AT:59:ARG:HH11	19:AT:59:ARG:CB	2.25	0.48
47:B0:28:SER:O	47:B0:29:VAL:HG12	2.13	0.48
47:B0:32:THR:HG22	47:B0:32:THR:O	2.13	0.48
23:BB:1315:C:H2'	23:BB:1316:U:H6	1.77	0.48
23:BB:1537:G:H2'	23:BB:1538:G:O4'	2.14	0.48
23:BB:1662:U:O2	23:BB:2687:U:H4'	2.13	0.48
23:BB:2526:G:H2'	23:BB:2527:C:H6	1.79	0.48
23:BB:2568:U:H2'	23:BB:2569:G:O4'	2.14	0.48
23:BB:2737:G:H2'	23:BB:2738:A:C8	2.48	0.48
23:BB:2845:U:H2'	23:BB:2846:G:C8	2.49	0.48
23:BB:388:G:N7	23:BB:390:U:H2'	2.28	0.48
23:BB:36:G:H4'	23:BB:451:U:C2	2.49	0.48
23:BB:503:A:H4'	23:BB:504:A:H5'	1.94	0.48
23:BB:921:C:H2'	23:BB:922:C:H6	1.79	0.48
23:BB:947:A:O2'	23:BB:984:A:H2	1.97	0.48
25:BC:79:ARG:NH2	25:BC:110:LYS:HD3	2.29	0.48
25:BC:64:VAL:HB	25:BC:65:ASP:H	1.42	0.48
26:BD:110:THR:HG22	26:BD:171:THR:CA	2.42	0.48
26:BD:38:LYS:N	26:BD:38:LYS:CD	2.76	0.48
28:BF:56:LEU:HA	28:BF:59:ILE:CG2	2.44	0.48
52:BI:37:PHE:CZ	52:BI:58:ILE:HD11	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:BJ:35:ARG:NH2	31:BJ:40:HIS:N	2.61	0.48
31:BJ:39:LYS:C	31:BJ:41:LYS:H	2.17	0.48
32:BK:85:VAL:O	32:BK:87:LEU:HD23	2.14	0.48
33:BL:27:LEU:N	33:BL:27:LEU:HD12	2.23	0.48
34:BM:7:THR:HG23	34:BM:91:TYR:OH	2.14	0.48
36:BO:40:ILE:CD1	36:BO:40:ILE:H	2.27	0.48
36:BO:46:GLU:C	36:BO:48:LEU:H	2.17	0.48
36:BO:87:ILE:CG1	36:BO:88:LYS:N	2.76	0.48
37:BP:76:HIS:CD2	37:BP:76:HIS:N	2.82	0.48
39:BR:1:MET:HA	39:BR:46:GLU:CB	2.43	0.48
23:BB:992:C:H1'	39:BR:91:GLN:CB	2.44	0.48
40:BS:6:LYS:CA	40:BS:104:THR:HA	2.32	0.48
40:BS:31:GLN:HB3	40:BS:35:ILE:CG2	2.44	0.48
23:BB:84:A:H3'	42:BU:5:ARG:HG3	1.96	0.48
24:BV:60:VAL:HG13	24:BV:72:VAL:O	2.13	0.48
46:BZ:3:LYS:CE	46:BZ:8:LYS:HA	2.41	0.48
1:CA:110:C:H2'	1:CA:111:G:O4'	2.14	0.48
1:CA:1343:G:H1'	8:CI:122:ARG:HH12	1.79	0.48
1:CA:1426:G:H2'	1:CA:1427:C:H6	1.79	0.48
1:CA:1480:A:H2'	1:CA:1481:U:H6	1.79	0.48
1:CA:627:G:H2'	1:CA:628:G:C8	2.48	0.48
1:CA:708:C:H2'	1:CA:709:U:H6	1.79	0.48
1:CA:737:C:H2'	1:CA:738:C:H6	1.79	0.48
1:CA:78:A:O2'	1:CA:79:G:H5'	2.14	0.48
20:CB:65:LYS:CB	20:CB:157:PRO:HA	2.43	0.48
3:CD:64:TYR:CE2	3:CD:93:LEU:HD13	2.48	0.48
9:CJ:40:ILE:CG1	9:CJ:73:LEU:HB3	2.44	0.48
22:DA:33:G:O2'	22:DA:34:A:H5'	2.14	0.48
23:DB:1165:A:H2'	23:DB:1166:G:H8	1.77	0.48
23:DB:1219:U:OP2	38:DQ:18:LYS:HE3	2.14	0.48
23:DB:137:U:H2'	23:DB:138:U:O4'	2.13	0.48
23:DB:1579:A:H2'	23:DB:1580:A:H8	1.78	0.48
23:DB:188:G:O2'	23:DB:189:G:H5'	2.14	0.48
23:DB:1903:G:OP1	25:DC:239:PHE:HA	2.13	0.48
23:DB:2350:C:O2'	23:DB:2351:G:H5'	2.14	0.48
23:DB:234:U:C2'	23:DB:235:U:H5'	2.44	0.48
23:DB:616:A:H3'	23:DB:617:G:C8	2.44	0.48
23:DB:942:G:H2'	23:DB:943:A:H8	1.79	0.48
25:DC:136:VAL:C	25:DC:165:ALA:HA	2.34	0.48
26:DD:105:LYS:O	26:DD:177:VAL:HG12	2.14	0.48
23:DB:607:U:P	27:DE:98:LYS:HG3	2.54	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DF:141:ASP:O	28:DF:142:TYR:HB3	2.13	0.48
30:DH:8:LYS:HE2	30:DH:9:VAL:N	2.27	0.48
31:DJ:67:ASN:C	31:DJ:69:ARG:H	2.16	0.48
32:DK:63:VAL:CG1	32:DK:103:VAL:HG12	2.44	0.48
33:DL:51:GLU:OE1	33:DL:51:GLU:N	2.47	0.48
33:DL:57:LEU:HB3	50:D3:54:LEU:HD22	1.95	0.48
33:DL:90:VAL:HG12	33:DL:122:VAL:HG22	1.96	0.48
23:DB:2376:A:H2	36:DO:117:PHE:HD2	1.61	0.48
37:DP:70:GLU:HG2	37:DP:71:ARG:HG2	1.96	0.48
1:AA:1356:G:H2'	1:AA:1357:A:C8	2.48	0.47
1:AA:434:U:H2'	1:AA:434:U:O2	2.15	0.47
1:AA:634:C:O2'	1:AA:635:A:H5'	2.14	0.47
1:AA:782:A:H2'	1:AA:783:C:O4'	2.14	0.47
20:AB:174:GLU:O	20:AB:178:LEU:HD23	2.14	0.47
20:AB:27:LYS:HB3	20:AB:28:PRO:HD3	1.96	0.47
6:AG:85:GLN:O	6:AG:86:VAL:C	2.52	0.47
12:AM:89:ARG:NH1	12:AM:94:LEU:HD13	2.29	0.47
51:B4:9:LYS:O	51:B4:10:LEU:HB3	2.14	0.47
23:BB:1299:G:H4'	23:BB:1301:A:C1'	2.42	0.47
23:BB:1681:G:N3	23:BB:1762:A:H2'	2.29	0.47
23:BB:1957:C:H2'	23:BB:1958:C:C6	2.48	0.47
23:BB:209:C:H2'	23:BB:210:C:C6	2.48	0.47
23:BB:2144:G:H2'	23:BB:2146:C:OP2	2.14	0.47
23:BB:2462:C:H2'	23:BB:2463:C:C6	2.49	0.47
23:BB:256:A:O2'	23:BB:257:C:H5'	2.14	0.47
23:BB:271:G:HO2'	23:BB:272:A:H8	1.62	0.47
23:BB:2800:A:H2'	23:BB:2801:G:C4'	2.43	0.47
23:BB:2815:C:H2'	23:BB:2816:G:H8	1.79	0.47
23:BB:2864:G:H2'	23:BB:2865:U:C6	2.49	0.47
23:BB:753:A:H2'	23:BB:754:U:H6	1.78	0.47
26:BD:16:THR:CG2	26:BD:17:GLU:H	2.14	0.47
26:BD:26:VAL:HG23	26:BD:187:LEU:HB3	1.96	0.47
26:BD:8:LYS:O	26:BD:9:VAL:HG22	2.13	0.47
28:BF:2:LYS:HB2	28:BF:100:GLU:HG2	1.96	0.47
29:BG:126:THR:CG2	29:BG:128:THR:HG23	2.44	0.47
29:BG:169:ARG:C	29:BG:171:LYS:H	2.17	0.47
29:BG:42:VAL:HG12	29:BG:51:PHE:CD1	2.49	0.47
31:BJ:125:TYR:HE2	31:BJ:134:ALA:HB2	1.78	0.47
23:BB:1132:U:H5'	31:BJ:85:LYS:HZ3	1.79	0.47
32:BK:63:VAL:CG1	32:BK:103:VAL:HA	2.44	0.47
34:BM:96:ILE:CG2	34:BM:97:GLN:H	2.17	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:118:ARG:NH1	47:B0:50:GLY:HA3	2.29	0.47
37:BP:81:ASP:O	37:BP:83:ILE:N	2.46	0.47
43:BW:40:ARG:HB2	43:BW:41:GLY:H	1.58	0.47
43:BW:44:PHE:CE2	43:BW:80:SER:HA	2.49	0.47
23:BB:852:U:H5'	45:BY:45:GLY:HA3	1.96	0.47
1:CA:1361:G:C2'	1:CA:1362:A:H5''	2.43	0.47
1:CA:836:G:H2'	1:CA:837:U:H6	1.79	0.47
1:CA:911:U:H2'	1:CA:912:C:C6	2.49	0.47
1:CA:921:U:H2'	1:CA:922:G:O4'	2.13	0.47
1:CA:954:G:H2'	1:CA:955:U:H6	1.79	0.47
20:CB:68:PHE:HA	20:CB:161:PHE:O	2.14	0.47
2:CC:146:LYS:HD3	2:CC:203:LYS:O	2.14	0.47
2:CC:140:ALA:HB2	2:CC:148:ILE:HD12	1.96	0.47
3:CD:118:SER:C	3:CD:120:LYS:H	2.18	0.47
3:CD:170:LEU:HD12	3:CD:170:LEU:O	2.14	0.47
8:CI:29:ILE:HD13	8:CI:37:TYR:CD2	2.49	0.47
14:CO:25:GLU:HA	14:CO:80:LEU:HD12	1.96	0.47
18:CS:43:MET:CB	18:CS:61:VAL:HG11	2.44	0.47
23:DB:2527:C:OP1	51:D4:34:LYS:HE2	2.14	0.47
23:DB:1531:C:H2'	23:DB:1532:A:H8	1.75	0.47
23:DB:1685:C:H2'	23:DB:1686:C:H6	1.78	0.47
23:DB:1681:G:N3	23:DB:1762:A:H2'	2.29	0.47
23:DB:1133:A:N6	23:DB:2025:C:O2'	2.46	0.47
23:DB:2143:C:H2'	23:DB:2144:G:C4'	2.43	0.47
23:DB:2220:U:H2'	23:DB:2221:G:C8	2.48	0.47
23:DB:2593:U:O2'	23:DB:2594:C:H5'	2.13	0.47
23:DB:2683:C:H2'	23:DB:2684:U:C6	2.49	0.47
23:DB:2795:C:O5'	23:DB:2795:C:H6	1.96	0.47
23:DB:396:G:H2'	23:DB:397:U:C6	2.49	0.47
23:DB:582:A:H2'	23:DB:583:G:C8	2.47	0.47
23:DB:704:G:O2'	23:DB:726:G:N2	2.46	0.47
23:DB:853:C:H2'	23:DB:854:C:H6	1.78	0.47
23:DB:871:U:H2'	23:DB:872:U:C6	2.49	0.47
23:DB:1658:C:OP1	26:DD:136:ASN:ND2	2.47	0.47
26:DD:148:GLN:HB3	26:DD:151:THR:HG21	1.96	0.47
26:DD:24:VAL:HG21	26:DD:193:VAL:CG1	2.44	0.47
28:DF:18:GLU:C	28:DF:20:ASN:N	2.66	0.47
30:DH:97:ARG:H	30:DH:97:ARG:HG2	1.38	0.47
32:DK:77:ILE:HD11	32:DK:105:ARG:HH12	1.77	0.47
32:DK:18:ARG:HB2	32:DK:45:GLU:CG	2.44	0.47
32:DK:63:VAL:HG22	32:DK:115:ILE:HD11	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:636:G:OP2	33:DL:126:ARG:NH2	2.47	0.47
33:DL:140:GLY:O	33:DL:142:ILE:N	2.47	0.47
33:DL:31:GLY:HA2	33:DL:36:LYS:HG2	1.96	0.47
36:DO:10:ARG:HG3	36:DO:96:GLY:H	1.77	0.47
36:DO:25:ARG:NE	36:DO:94:ARG:HH12	2.06	0.47
44:DX:51:ALA:O	44:DX:53:VAL:HG12	2.14	0.47
45:DY:13:ILE:HD13	45:DY:13:ILE:H	1.79	0.47
1:AA:1126:U:O2'	1:AA:1127:G:H5'	2.14	0.47
1:AA:1198:G:H2'	1:AA:1199:U:C6	2.49	0.47
1:AA:1221:G:OP1	1:AA:1320:C:N4	2.47	0.47
1:AA:152:A:N6	1:AA:170:U:C2	2.83	0.47
1:AA:197:A:H4'	1:AA:198:G:O5'	2.15	0.47
1:AA:470:C:H2'	1:AA:471:U:H6	1.79	0.47
1:AA:737:C:H2'	1:AA:738:C:H6	1.79	0.47
20:AB:128:LEU:CG	20:AB:132:GLU:HG2	2.44	0.47
20:AB:128:LEU:HD13	20:AB:129:THR:H	1.79	0.47
20:AB:116:LEU:HB3	20:AB:140:LEU:CD2	2.44	0.47
20:AB:209:VAL:HG23	20:AB:210:THR:H	1.77	0.47
3:AD:67:LEU:N	3:AD:67:LEU:HD22	2.29	0.47
4:AE:132:PRO:HA	4:AE:135:VAL:HG22	1.95	0.47
4:AE:134:ASN:O	4:AE:137:ARG:HG2	2.13	0.47
5:AF:18:VAL:HG11	5:AF:58:HIS:CE1	2.49	0.47
5:AF:35:LYS:HD2	5:AF:65:GLU:OE2	2.14	0.47
6:AG:102:TRP:O	6:AG:106:ALA:HB3	2.14	0.47
8:AI:24:ASN:CA	8:AI:26:LYS:HZ2	2.27	0.47
8:AI:59:LYS:HB3	8:AI:59:LYS:HZ3	1.78	0.47
1:AA:1225:A:H5'	12:AM:101:THR:HG23	1.96	0.47
13:AN:27:LYS:HE2	13:AN:48:GLN:NE2	2.28	0.47
23:BB:1358:G:N2	23:BB:1372:U:C5	2.82	0.47
23:BB:1427:A:H4'	23:BB:1428:C:O4'	2.14	0.47
23:BB:1585:C:H2'	23:BB:1586:A:O4'	2.14	0.47
23:BB:2087:G:H2'	23:BB:2088:A:C8	2.49	0.47
23:BB:2144:G:C4	23:BB:2146:C:H5'	2.49	0.47
23:BB:233:A:H61	23:BB:428:A:N6	2.12	0.47
23:BB:284:U:H2'	23:BB:285:G:C8	2.48	0.47
23:BB:863:A:H2'	23:BB:864:G:C8	2.49	0.47
23:BB:946:C:H2'	23:BB:947:A:C8	2.49	0.47
25:BC:131:MET:O	25:BC:134:ILE:HB	2.14	0.47
25:BC:141:HIS:O	25:BC:153:LEU:HD21	2.14	0.47
26:BD:206:ALA:C	26:BD:208:LYS:H	2.17	0.47
28:BF:177:ARG:HG2	28:BF:178:LYS:N	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:33:ILE:HG22	28:BF:155:ILE:HG12	1.95	0.47
29:BG:10:VAL:HG13	29:BG:10:VAL:O	2.13	0.47
29:BG:5:LYS:NZ	29:BG:69:ALA:HB2	2.29	0.47
30:BH:141:LYS:N	30:BH:141:LYS:HD3	2.29	0.47
52:BI:91:LYS:O	52:BI:91:LYS:HG3	2.13	0.47
31:BJ:7:LYS:C	31:BJ:9:GLU:H	2.17	0.47
33:BL:126:ARG:HD3	33:BL:126:ARG:C	2.34	0.47
33:BL:39:LYS:HB2	33:BL:46:VAL:HG22	1.95	0.47
33:BL:90:VAL:HG12	33:BL:121:THR:O	2.14	0.47
34:BM:33:LEU:HB2	34:BM:101:VAL:HG22	1.96	0.47
34:BM:80:VAL:O	34:BM:81:ARG:HB3	2.14	0.47
37:BP:112:ARG:H	37:BP:112:ARG:HH11	1.60	0.47
39:BR:18:GLN:N	39:BR:18:GLN:OE1	2.47	0.47
41:BT:60:THR:HG22	41:BT:83:ALA:CB	2.44	0.47
41:BT:93:LEU:H	41:BT:93:LEU:CD2	2.25	0.47
43:BW:46:ALA:HA	43:BW:57:THR:OG1	2.14	0.47
1:CA:1302:C:H4'	1:CA:1303:C:OP1	2.13	0.47
1:CA:1404:C:H2'	1:CA:1405:G:H8	1.80	0.47
1:CA:370:C:H2'	1:CA:371:A:H8	1.77	0.47
1:CA:413:G:H5'	1:CA:414:A:OP1	2.14	0.47
1:CA:587:G:H4'	7:CH:3:GLN:HA	1.96	0.47
20:CB:80:LYS:HA	20:CB:90:PHE:CZ	2.50	0.47
2:CC:23:ALA:HB3	2:CC:28:PHE:HD1	1.78	0.47
2:CC:2:GLN:N	2:CC:2:GLN:NE2	2.62	0.47
2:CC:71:ARG:HG3	2:CC:71:ARG:O	2.15	0.47
5:CF:51:ILE:O	5:CF:51:ILE:HG23	2.14	0.47
5:CF:6:ILE:HG23	5:CF:62:MET:CB	2.35	0.47
10:CK:12:ARG:HB3	10:CK:13:LYS:HD2	1.96	0.47
1:CA:500:G:C5'	11:CL:120:ARG:HH21	2.28	0.47
11:CL:80:LEU:HB3	11:CL:97:VAL:CG2	2.44	0.47
14:CO:81:ILE:HD12	14:CO:87:ARG:HB2	1.96	0.47
49:D2:12:ARG:NH2	49:D2:16:HIS:HB2	2.28	0.47
22:DA:47:C:H5'	36:DO:97:PHE:HZ	1.77	0.47
23:DB:1213:A:N6	23:DB:1236:G:H1'	2.29	0.47
23:DB:1358:G:N2	23:DB:1372:U:C5	2.82	0.47
23:DB:137:U:H2'	23:DB:138:U:C5'	2.45	0.47
23:DB:1433:A:H2'	23:DB:1434:A:O4'	2.13	0.47
23:DB:1585:C:H2'	23:DB:1586:A:O4'	2.14	0.47
23:DB:16:C:O3'	47:D0:10:SER:HA	2.14	0.47
23:DB:1775:U:H2'	23:DB:1776:G:O4'	2.14	0.47
1:CA:1517:G:H1'	23:DB:1919:A:O3'	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2207:C:H2'	23:DB:2208:C:C6	2.48	0.47
23:DB:2526:G:H2'	23:DB:2527:C:H6	1.79	0.47
23:DB:2519:U:H5'	23:DB:2567:G:N2	2.29	0.47
23:DB:2647:U:O2'	23:DB:2648:G:H5'	2.15	0.47
23:DB:2730:C:H4'	26:DD:174:SER:O	2.13	0.47
23:DB:2828:G:O2'	23:DB:2829:A:H5'	2.14	0.47
23:DB:298:G:C2	23:DB:339:U:C5	3.02	0.47
23:DB:356:G:H2'	23:DB:357:C:C6	2.50	0.47
23:DB:460:A:H2'	23:DB:461:C:O4'	2.14	0.47
23:DB:491:G:H2'	23:DB:492:A:O4'	2.14	0.47
23:DB:543:G:C5	23:DB:544:C:H1'	2.49	0.47
23:DB:544:C:O5'	23:DB:545:U:OP1	2.32	0.47
23:DB:664:G:O2'	23:DB:665:U:H5'	2.14	0.47
23:DB:806:C:H2'	23:DB:807:U:C6	2.49	0.47
23:DB:844:A:C2	23:DB:845:A:C6	3.02	0.47
25:DC:123:ILE:HG12	25:DC:135:PRO:HD2	1.97	0.47
26:DD:180:VAL:HG23	26:DD:181:ASP:H	1.78	0.47
26:DD:95:SER:O	26:DD:96:ILE:HG13	2.12	0.47
29:DG:10:VAL:HG13	29:DG:14:VAL:HG12	1.96	0.47
29:DG:10:VAL:HG13	29:DG:14:VAL:CG1	2.44	0.47
30:DH:135:HIS:HB3	30:DH:138:VAL:CB	2.44	0.47
31:DJ:133:ALA:O	31:DJ:135:GLN:N	2.47	0.47
23:DB:956:G:H4'	34:DM:81:ARG:HH21	1.79	0.47
35:DN:2:ARG:NH2	35:DN:4:ARG:HH11	2.12	0.47
35:DN:36:THR:HG22	35:DN:37:THR:H	1.78	0.47
37:DP:79:VAL:HG23	37:DP:81:ASP:OD1	2.14	0.47
38:DQ:116:LEU:HD22	38:DQ:116:LEU:N	2.28	0.47
41:DT:31:VAL:HG23	41:DT:82:LYS:HG3	1.95	0.47
41:DT:54:GLU:HB3	41:DT:91:GLN:OE1	2.14	0.47
1:AA:528:C:H41	11:AL:45:ASN:CG	2.17	0.47
1:AA:790:A:N1	1:AA:1497:G:H5''	2.29	0.47
2:AC:171:ARG:NH1	2:AC:173:PRO:HG3	2.29	0.47
2:AC:71:ARG:O	2:AC:74:ILE:HG22	2.13	0.47
3:AD:81:LEU:CD1	3:AD:92:LEU:HD21	2.44	0.47
7:AH:34:ALA:HB1	7:AH:109:VAL:HG21	1.95	0.47
9:AJ:48:ARG:HG2	9:AJ:48:ARG:HH11	1.77	0.47
9:AJ:5:ARG:HB3	9:AJ:77:VAL:HA	1.96	0.47
10:AK:17:ASP:HA	10:AK:80:ASN:O	2.14	0.47
15:AP:67:ILE:HD11	15:AP:72:ALA:HA	1.95	0.47
16:AQ:10:ARG:NE	16:AQ:11:VAL:H	2.12	0.47
16:AQ:43:LEU:HD21	16:AQ:72:TRP:CE2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:B1:31:GLU:H	48:B1:32:LYS:NZ	2.12	0.47
23:BB:1438:U:O2'	23:BB:1439:A:H5'	2.14	0.47
23:BB:1633:G:O2'	23:BB:1634:A:H5''	2.14	0.47
23:BB:2350:C:O2'	23:BB:2351:G:H5'	2.13	0.47
23:BB:2822:G:O6	35:BN:1:MET:N	2.41	0.47
23:BB:2893:A:H4'	23:BB:2894:G:H5'	1.96	0.47
23:BB:346:A:H5'	23:BB:346:A:N3	2.30	0.47
23:BB:835:C:H5'	50:B3:50:SER:HB3	1.95	0.47
25:BC:24:HIS:ND1	25:BC:25:LYS:N	2.62	0.47
25:BC:29:PHE:C	25:BC:29:PHE:CD1	2.87	0.47
25:BC:63:ILE:HG22	25:BC:64:VAL:N	2.29	0.47
27:BE:165:HIS:C	27:BE:166:LYS:HG3	2.35	0.47
27:BE:147:LEU:HA	27:BE:166:LYS:O	2.13	0.47
27:BE:73:ILE:C	27:BE:75:SER:N	2.66	0.47
55:BB:3339:HOH:O	27:BE:81:GLY:HA2	2.14	0.47
28:BF:39:VAL:O	28:BF:39:VAL:HG12	2.14	0.47
29:BG:42:VAL:HG12	29:BG:51:PHE:HD1	1.78	0.47
29:BG:32:LEU:HD21	29:BG:74:MET:SD	2.54	0.47
30:BH:6:LEU:CD2	30:BH:50:ARG:HD3	2.44	0.47
30:BH:5:LEU:C	30:BH:7:ASP:H	2.18	0.47
52:BI:116:MET:HE2	52:BI:124:MET:HA	1.97	0.47
22:BA:103:U:H4'	34:BM:136:MET:HE2	1.96	0.47
35:BN:75:ILE:C	35:BN:75:ILE:HD13	2.35	0.47
23:BB:812:C:H4'	38:BQ:12:ARG:HH22	1.79	0.47
41:BT:33:LYS:HE3	41:BT:80:TRP:NE1	2.29	0.47
1:CA:340:U:H2'	1:CA:341:C:H6	1.78	0.47
1:CA:113:G:H1'	1:CA:354:G:H5'	1.96	0.47
1:CA:470:C:H2'	1:CA:471:U:H6	1.79	0.47
1:CA:781:A:H2'	1:CA:782:A:C5'	2.38	0.47
1:CA:803:G:H2'	1:CA:804:U:C6	2.49	0.47
1:CA:893:C:H2'	1:CA:894:G:C8	2.49	0.47
20:CB:95:TRP:CH2	20:CB:99:MET:HB2	2.49	0.47
2:CC:123:LEU:HD23	2:CC:195:ILE:HD12	1.96	0.47
4:CE:89:THR:HG23	4:CE:90:GLY:N	2.29	0.47
6:CG:94:ARG:HG2	6:CG:98:LEU:HD11	1.95	0.47
11:CL:82:ARG:HG2	11:CL:82:ARG:HH11	1.79	0.47
16:CQ:58:VAL:HB	16:CQ:74:LEU:CD2	2.43	0.47
18:CS:39:ILE:CG2	18:CS:66:VAL:HA	2.45	0.47
18:CS:18:VAL:C	18:CS:42:ASN:HD21	2.17	0.47
19:CT:54:GLN:N	19:CT:55:PRO:HD2	2.30	0.47
1:CA:1527:U:OP2	21:CU:38:GLU:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1027:A:H2	23:DB:2488:G:H4'	1.79	0.47
23:DB:1515:A:H5'	23:DB:1557:C:C5'	2.44	0.47
23:DB:1567:G:H5''	25:DC:84:PRO:HG3	1.96	0.47
23:DB:1599:U:H2'	23:DB:1600:C:C6	2.48	0.47
23:DB:1957:C:H2'	23:DB:1958:C:C6	2.49	0.47
23:DB:2286:G:H5'	23:DB:2286:G:C8	2.49	0.47
23:DB:263:G:H2'	23:DB:264:C:O4'	2.14	0.47
23:DB:2845:U:H2'	23:DB:2846:G:C8	2.49	0.47
23:DB:851:C:O2'	23:DB:852:U:H5'	2.14	0.47
25:DC:162:GLN:HE21	25:DC:174:ARG:NH2	2.12	0.47
25:DC:12:ARG:HB2	25:DC:20:ASN:HA	1.96	0.47
25:DC:266:ILE:HD11	25:DC:268:ARG:HH22	1.79	0.47
25:DC:42:ARG:HE	25:DC:42:ARG:CA	2.27	0.47
26:DD:11:MET:O	26:DD:23:PRO:HD2	2.15	0.47
28:DF:163:GLU:HG2	28:DF:166:ARG:NH1	2.28	0.47
30:DH:3:VAL:CG1	30:DH:37:VAL:HG11	2.44	0.47
30:DH:51:ARG:HG2	30:DH:51:ARG:NH1	2.27	0.47
23:DB:1666:G:OP1	32:DK:66:LYS:HD2	2.14	0.47
33:DL:39:LYS:NZ	33:DL:39:LYS:CA	2.76	0.47
34:DM:57:VAL:CG1	34:DM:58:LYS:H	2.12	0.47
1:AA:1469:C:H2'	1:AA:1470:U:O4'	2.15	0.47
1:AA:1504:G:H4'	1:AA:1505:G:N9	2.29	0.47
1:AA:878:A:H1'	7:AH:3:GLN:HE22	1.79	0.47
20:AB:113:LEU:O	20:AB:113:LEU:HD23	2.15	0.47
20:AB:95:TRP:CZ3	20:AB:171:ALA:HA	2.49	0.47
2:AC:111:ASP:OD2	2:AC:114:LEU:HG	2.14	0.47
2:AC:141:MET:HE1	2:AC:147:GLY:N	2.18	0.47
2:AC:40:GLN:HE22	2:AC:44:LYS:HD2	1.80	0.47
4:AE:150:GLU:C	4:AE:152:VAL:H	2.17	0.47
4:AE:40:ASP:C	4:AE:42:ASN:H	2.17	0.47
5:AF:63:ASN:HD22	5:AF:96:VAL:HB	1.79	0.47
12:AM:109:LYS:HG3	12:AM:110:GLY:H	1.78	0.47
12:AM:89:ARG:HH11	12:AM:94:LEU:HD13	1.79	0.47
13:AN:84:ARG:HH11	13:AN:84:ARG:HG3	1.80	0.47
15:AP:42:ILE:CB	15:AP:46:LYS:HD2	2.29	0.47
18:AS:40:PHE:HB3	18:AS:43:MET:HG3	1.96	0.47
23:BB:2615:U:C4	47:B0:2:VAL:O	2.68	0.47
23:BB:1239:G:H2'	23:BB:1240:U:C6	2.50	0.47
23:BB:1328:A:H2'	23:BB:1330:C:C4	2.50	0.47
23:BB:160:A:H2'	23:BB:161:A:C8	2.49	0.47
23:BB:1847:A:H4'	23:BB:1848:A:H8	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2222:C:O2'	23:BB:2223:G:H5'	2.14	0.47
23:BB:2344:U:H4'	23:BB:2345:G:OP1	2.14	0.47
23:BB:2393:U:O2'	23:BB:2394:C:H5'	2.15	0.47
23:BB:2647:U:O2'	23:BB:2648:G:H5'	2.14	0.47
23:BB:2757:A:N3	23:BB:2757:A:H2'	2.28	0.47
23:BB:338:G:N2	23:BB:339:U:H1'	2.29	0.47
23:BB:640:C:H1'	50:B3:44:ARG:HH21	1.80	0.47
23:BB:702:U:H2'	23:BB:703:U:C6	2.49	0.47
23:BB:806:C:H2'	23:BB:807:U:C6	2.50	0.47
25:BC:66:PHE:O	25:BC:68:ARG:N	2.47	0.47
26:BD:119:ALA:HA	26:BD:123:LYS:HZ3	1.78	0.47
26:BD:129:THR:O	26:BD:130:GLN:CB	2.62	0.47
23:BB:2572:A:OP1	26:BD:151:THR:CG2	2.62	0.47
26:BD:61:THR:H	26:BD:62:LYS:HE3	1.78	0.47
27:BE:118:LEU:HD11	27:BE:187:VAL:HG12	1.96	0.47
27:BE:159:LEU:O	27:BE:159:LEU:HD23	2.15	0.47
23:BB:675:A:OP1	27:BE:58:LYS:HG2	2.14	0.47
30:BH:132:PHE:CB	30:BH:140:ALA:HB3	2.44	0.47
31:BJ:113:PRO:O	31:BJ:115:GLY:N	2.44	0.47
31:BJ:46:PRO:O	31:BJ:48:VAL:N	2.43	0.47
23:BB:1131:G:H1'	31:BJ:85:LYS:NZ	2.28	0.47
33:BL:25:SER:HB3	33:BL:27:LEU:HG	1.97	0.47
33:BL:55:MET:N	33:BL:56:PRO:CD	2.78	0.47
33:BL:61:LEU:O	33:BL:62:PRO:O	2.32	0.47
35:BN:106:ASP:O	35:BN:108:ALA:N	2.47	0.47
35:BN:44:LEU:HD23	35:BN:113:ILE:CG1	2.37	0.47
41:BT:40:LYS:HE3	41:BT:59:ASN:HA	1.96	0.47
41:BT:86:THR:C	41:BT:87:LEU:HD22	2.35	0.47
42:BU:35:VAL:C	42:BU:36:GLU:HG2	2.34	0.47
23:BB:399:U:OP2	46:BZ:49:ARG:NH1	2.48	0.47
1:CA:1524:C:H2'	1:CA:1525:G:H8	1.78	0.47
1:CA:36:C:O2'	1:CA:37:U:H5'	2.15	0.47
1:CA:719:C:O2'	17:CR:37:LYS:HD3	2.15	0.47
1:CA:954:G:H2'	1:CA:955:U:C6	2.50	0.47
2:CC:155:ARG:NH2	2:CC:160:GLU:HA	2.29	0.47
3:CD:15:GLY:HA2	3:CD:34:GLU:HB2	1.97	0.47
3:CD:171:GLU:HG3	3:CD:182:LYS:CG	2.45	0.47
4:CE:114:LEU:HB3	4:CE:119:VAL:HB	1.96	0.47
8:CI:7:GLY:HA3	8:CI:81:GLY:O	2.14	0.47
9:CJ:40:ILE:HD11	9:CJ:73:LEU:HB3	1.97	0.47
1:CA:1060:U:OP1	9:CJ:53:ILE:HD11	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:51:VAL:O	9:CJ:62:ARG:HA	2.14	0.47
14:CO:26:VAL:HG12	14:CO:30:LEU:HD12	1.96	0.47
16:CQ:13:SER:CB	16:CQ:21:VAL:HB	2.41	0.47
16:CQ:8:GLN:HA	16:CQ:59:GLU:HA	1.96	0.47
17:CR:25:ILE:HD12	17:CR:25:ILE:O	2.15	0.47
47:D0:41:HIS:CE1	47:D0:42:ILE:HG22	2.49	0.47
23:DB:52:A:C5	23:DB:118:A:C2	3.02	0.47
23:DB:1561:C:H2'	23:DB:1562:U:C6	2.49	0.47
23:DB:1814:G:C5'	25:DC:51:ARG:HG2	2.44	0.47
23:DB:2383:G:N7	50:D3:35:LYS:HD3	2.30	0.47
23:DB:2868:A:H2'	23:DB:2869:G:H8	1.78	0.47
23:DB:545:U:H2'	23:DB:547:A:OP1	2.15	0.47
23:DB:825:A:H2'	23:DB:826:U:O4'	2.14	0.47
23:DB:832:U:H2'	23:DB:833:A:C8	2.49	0.47
25:DC:61:TYR:CE1	25:DC:63:ILE:HD11	2.49	0.47
26:DD:116:LYS:N	26:DD:116:LYS:HD2	2.28	0.47
23:DB:2052:A:H5'	26:DD:146:ILE:O	2.14	0.47
26:DD:35:THR:HB	26:DD:48:ILE:CG1	2.44	0.47
27:DE:48:THR:HG23	27:DE:85:PHE:N	2.29	0.47
30:DH:42:LYS:O	30:DH:45:GLU:HB3	2.14	0.47
30:DH:6:LEU:HB2	30:DH:35:LYS:CB	2.43	0.47
31:DJ:14:ASP:HB3	31:DJ:16:TYR:HD1	1.79	0.47
32:DK:19:VAL:C	32:DK:41:ILE:HD11	2.34	0.47
32:DK:7:MET:SD	32:DK:20:MET:HB2	2.55	0.47
37:DP:100:ARG:HH11	37:DP:100:ARG:CB	2.17	0.47
37:DP:27:VAL:HG21	37:DP:84:SER:O	2.14	0.47
39:DR:18:GLN:CA	39:DR:99:THR:HA	2.44	0.47
35:DN:107:ASN:ND2	40:DS:40:ASN:HD22	2.10	0.47
36:DO:15:ARG:HH12	43:DW:76:ARG:CD	2.27	0.47
45:DY:50:VAL:HG12	45:DY:53:MET:HG2	1.96	0.47
46:DZ:49:ARG:CZ	46:DZ:49:ARG:HB2	2.45	0.47
1:AA:1002:G:N3	1:AA:1002:G:H2'	2.29	0.47
1:AA:1014:A:C2	1:AA:1219:A:H1'	2.50	0.47
1:AA:1257:A:H2'	1:AA:1257:A:N3	2.28	0.47
1:AA:49:U:O2'	1:AA:50:A:H2'	2.15	0.47
1:AA:699:C:C3'	1:AA:700:G:H5''	2.44	0.47
1:AA:831:A:O2'	1:AA:832:G:H5'	2.15	0.47
1:AA:985:C:H2'	1:AA:986:U:C6	2.49	0.47
3:AD:116:LEU:HB3	3:AD:122:ILE:HD11	1.97	0.47
4:AE:84:VAL:HG22	4:AE:85:LYS:N	2.28	0.47
6:AG:19:SER:HB2	6:AG:21:LEU:HD21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AJ:37:ARG:HB2	9:AJ:77:VAL:HG11	1.97	0.47
13:AN:17:ASP:HA	13:AN:21:ALA:HB3	1.95	0.47
18:AS:51:HIS:HB2	18:AS:56:HIS:CE1	2.50	0.47
48:B1:15:GLY:CA	48:B1:47:ILE:HD12	2.44	0.47
48:B1:47:ILE:HG13	48:B1:48:TYR:H	1.77	0.47
22:BA:15:A:H1'	22:BA:109:A:N7	2.30	0.47
23:BB:1175:A:H2'	23:BB:1176:U:H5'	1.95	0.47
23:BB:1827:U:O2'	23:BB:1828:G:H5'	2.14	0.47
23:BB:2331:G:H21	23:BB:2336:A:H8	1.62	0.47
23:BB:2828:G:O2'	23:BB:2829:A:H5'	2.14	0.47
23:BB:2845:U:H5''	37:BP:54:LEU:HA	1.96	0.47
23:BB:2880:C:O2	35:BN:93:GLY:HA3	2.14	0.47
23:BB:671:C:OP1	33:BL:48:ARG:NH1	2.48	0.47
23:BB:716:A:C3'	23:BB:717:C:H5''	2.44	0.47
25:BC:155:ARG:HB3	25:BC:159:THR:HG21	1.95	0.47
25:BC:22:GLU:HA	25:BC:202:ARG:NH1	2.29	0.47
26:BD:37:VAL:HG22	26:BD:46:ARG:CD	2.44	0.47
27:BE:73:ILE:O	27:BE:73:ILE:HG22	2.15	0.47
28:BF:93:GLU:OE1	28:BF:94:ARG:N	2.47	0.47
30:BH:96:THR:HB	30:BH:112:LYS:HB2	1.97	0.47
52:BI:48:ILE:O	52:BI:49:GLU:HB3	2.14	0.47
33:BL:62:PRO:HB3	50:B3:24:LYS:HE2	1.95	0.47
34:BM:116:ALA:HA	34:BM:119:LEU:HD23	1.97	0.47
35:BN:33:ILE:HG22	35:BN:113:ILE:O	2.14	0.47
37:BP:106:ALA:O	37:BP:108:ARG:HD2	2.14	0.47
38:BQ:33:VAL:O	38:BQ:34:ALA:C	2.53	0.47
40:BS:71:VAL:HG22	40:BS:107:VAL:CG1	2.32	0.47
23:BB:2014:A:H4'	40:BS:93:ALA:CB	2.45	0.47
41:BT:30:ILE:CG2	41:BT:32:LEU:HD13	2.44	0.47
42:BU:11:ILE:HG22	42:BU:12:VAL:N	2.30	0.47
24:BV:53:LYS:CE	24:BV:55:GLU:HG3	2.45	0.47
43:BW:22:VAL:HG12	43:BW:23:LYS:H	1.79	0.47
45:BY:54:VAL:HG22	45:BY:55:LYS:N	2.24	0.47
46:BZ:30:HIS:HB2	46:BZ:48:GLN:HB3	1.96	0.47
46:BZ:50:ASP:HA	46:BZ:53:THR:CG2	2.44	0.47
1:CA:1086:U:H3	1:CA:1099:G:N2	2.03	0.47
1:CA:333:U:H2'	1:CA:334:C:C6	2.49	0.47
1:CA:470:C:O2'	1:CA:471:U:H5'	2.14	0.47
1:CA:563:A:H2'	1:CA:567:G:C8	2.50	0.47
1:CA:946:A:H2'	1:CA:947:G:H8	1.73	0.47
1:CA:987:G:H2'	1:CA:988:G:H8	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:CB:172:ILE:CD1	20:CB:182:VAL:HG11	2.45	0.47
20:CB:80:LYS:HA	20:CB:90:PHE:CE1	2.48	0.47
2:CC:52:SER:HB3	2:CC:114:LEU:HD11	1.95	0.47
2:CC:37:LYS:O	2:CC:40:GLN:HG2	2.14	0.47
4:CE:67:ARG:HH11	4:CE:67:ARG:CB	2.26	0.47
8:CI:74:GLN:O	8:CI:78:ILE:HG13	2.14	0.47
11:CL:86:VAL:CG1	11:CL:89:LEU:HD23	2.44	0.47
15:CP:52:LEU:HD11	15:CP:74:LEU:HD13	1.95	0.47
21:CU:48:LYS:C	21:CU:48:LYS:HD2	2.35	0.47
21:CU:4:LYS:O	21:CU:5:VAL:HG23	2.14	0.47
21:CU:8:ASN:HB2	21:CU:9:GLU:CD	2.34	0.47
47:D0:36:LYS:HG3	47:D0:40:HIS:O	2.14	0.47
49:D2:18:PHE:O	49:D2:22:MET:N	2.45	0.47
23:DB:687:C:H1'	49:D2:4:THR:HG22	1.97	0.47
22:DA:23:G:N2	22:DA:24:G:H1	2.12	0.47
23:DB:1109:C:O2'	23:DB:1110:G:H5'	2.15	0.47
23:DB:137:U:H2'	23:DB:138:U:H5'	1.96	0.47
23:DB:1633:G:O2'	23:DB:1634:A:H5''	2.14	0.47
23:DB:1657:U:OP1	26:DD:141:ARG:HB2	2.14	0.47
23:DB:1818:U:H5''	25:DC:155:ARG:CG	2.44	0.47
23:DB:2630:G:H2'	23:DB:2631:G:C8	2.50	0.47
23:DB:2815:C:H2'	23:DB:2816:G:H8	1.79	0.47
23:DB:2834:G:H2'	23:DB:2879:A:H61	1.78	0.47
23:DB:82:U:H5''	23:DB:296:U:H5''	1.96	0.47
25:DC:86:ARG:C	25:DC:155:ARG:NH1	2.68	0.47
25:DC:224:MET:HG3	25:DC:233:GLY:CA	2.45	0.47
26:DD:129:THR:O	26:DD:130:GLN:CB	2.62	0.47
23:DB:2680:U:C5'	26:DD:194:PRO:HA	2.44	0.47
27:DE:4:VAL:HA	27:DE:14:VAL:CG2	2.45	0.47
28:DF:36:ASN:ND2	28:DF:87:LYS:H	2.12	0.47
28:DF:41:GLU:H	28:DF:45:ASP:HB3	1.79	0.47
52:DI:17:ALA:O	52:DI:18:ASN:HB3	2.14	0.47
31:DJ:25:LEU:H	31:DJ:25:LEU:HD22	1.79	0.47
33:DL:118:THR:C	33:DL:120:VAL:N	2.64	0.47
34:DM:133:LYS:HD2	34:DM:134:THR:N	2.29	0.47
34:DM:96:ILE:HG22	34:DM:97:GLN:HE22	1.78	0.47
34:DM:8:LYS:HG2	34:DM:9:PHE:N	2.28	0.47
39:DR:47:VAL:HG13	39:DR:49:ILE:H	1.80	0.47
39:DR:79:ARG:CD	39:DR:81:LYS:HG3	2.44	0.47
40:DS:72:THR:HG23	40:DS:73:LYS:N	2.29	0.47
23:DB:2270:A:H4'	43:DW:18:LYS:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DW:66:VAL:HG13	43:DW:67:LYS:N	2.24	0.47
44:DX:18:LEU:CD2	44:DX:18:LEU:H	2.14	0.47
46:DZ:33:ASN:HB3	46:DZ:46:GLY:HA2	1.96	0.47
1:AA:1072:G:H2'	1:AA:1073:U:C6	2.49	0.47
1:AA:1110:A:C2'	1:AA:1111:A:H5'	2.44	0.47
1:AA:1250:A:H4'	8:AI:69:GLY:N	2.24	0.47
1:AA:1409:C:N4	1:AA:1491:G:N1	2.50	0.47
1:AA:472:U:N3	1:AA:473:U:C4	2.82	0.47
1:AA:86:G:H4'	1:AA:86:G:OP1	2.13	0.47
1:AA:947:G:H2'	1:AA:948:C:C6	2.49	0.47
20:AB:113:LEU:HD12	20:AB:143:LEU:HB3	1.97	0.47
20:AB:162:VAL:O	20:AB:184:ALA:HA	2.15	0.47
6:AG:112:ASP:CB	6:AG:118:ARG:HG2	2.39	0.47
6:AG:47:GLU:OE1	6:AG:47:GLU:HA	2.13	0.47
7:AH:111:THR:HG23	7:AH:114:ALA:CB	2.45	0.47
9:AJ:39:PRO:HB3	9:AJ:74:VAL:HG22	1.96	0.47
11:AL:24:GLU:HB3	11:AL:26:CYS:SG	2.54	0.47
15:AP:33:ILE:HD12	15:AP:33:ILE:H	1.80	0.47
15:AP:76:LYS:HD2	15:AP:80:LYS:HD2	1.97	0.47
50:B3:24:LYS:CE	50:B3:29:ARG:HH22	2.27	0.47
22:BA:58:A:H2'	22:BA:59:A:O4'	2.14	0.47
23:BB:1188:U:O2'	23:BB:1189:A:H5'	2.14	0.47
23:BB:142:A:O2'	41:BT:2:ILE:HG21	2.14	0.47
23:BB:1882:U:O2'	23:BB:1883:U:H5'	2.14	0.47
23:BB:2205:A:H2'	23:BB:2206:C:C6	2.49	0.47
23:BB:299:A:N6	23:BB:322:A:H1'	2.30	0.47
23:BB:423:A:H5'	23:BB:424:G:H5'	1.96	0.47
23:BB:832:U:H2'	23:BB:833:A:C8	2.49	0.47
23:BB:859:G:O2'	23:BB:860:U:OP2	2.33	0.47
26:BD:41:ALA:O	26:BD:42:ASN:HB3	2.15	0.47
28:BF:16:MET:O	28:BF:20:ASN:HA	2.14	0.47
22:BA:42:C:H4'	28:BF:63:LYS:O	2.13	0.47
32:BK:64:ARG:HB3	32:BK:79:PHE:HB2	1.97	0.47
23:BB:956:G:C1'	34:BM:81:ARG:HH22	2.27	0.47
37:BP:94:ALA:HB3	37:BP:99:LEU:HD13	1.96	0.47
41:BT:64:LYS:HZ3	41:BT:79:ASP:HB2	1.80	0.47
41:BT:63:VAL:CG1	41:BT:80:TRP:HB2	2.45	0.47
43:BW:17:ALA:O	43:BW:35:ILE:HA	2.15	0.47
1:CA:1250:A:H2'	1:CA:1251:A:H8	1.79	0.47
1:CA:1301:U:O2'	1:CA:1302:C:P	2.72	0.47
1:CA:152:A:N6	1:CA:170:U:C2	2.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:202:G:H1'	1:CA:468:A:H8	1.79	0.47
1:CA:512:U:H2'	1:CA:513:C:C6	2.49	0.47
1:CA:589:U:O2'	1:CA:590:U:H5'	2.14	0.47
1:CA:792:A:H1'	1:CA:794:A:N7	2.29	0.47
20:CB:96:LEU:HB2	20:CB:99:MET:HG3	1.96	0.47
6:CG:55:LYS:HB3	6:CG:59:GLU:CD	2.35	0.47
7:CH:50:VAL:HG13	7:CH:50:VAL:O	2.15	0.47
10:CK:70:ALA:HB1	10:CK:74:LYS:CD	2.41	0.47
11:CL:82:ARG:NH2	11:CL:95:HIS:ND1	2.63	0.47
12:CM:47:LEU:HD23	12:CM:47:LEU:N	2.29	0.47
15:CP:61:VAL:CA	15:CP:65:ALA:HB3	2.44	0.47
19:CT:31:ILE:HD13	19:CT:74:HIS:CE1	2.49	0.47
47:D0:12:ARG:HH21	47:D0:16:ARG:HG3	1.80	0.47
47:D0:24:VAL:O	47:D0:24:VAL:HG12	2.15	0.47
47:D0:29:VAL:HB	47:D0:34:GLY:CA	2.44	0.47
22:DA:64:G:H2'	22:DA:65:U:C6	2.49	0.47
23:DB:159:G:O2'	23:DB:160:A:H5''	2.15	0.47
23:DB:2331:G:H21	23:DB:2336:A:H8	1.62	0.47
23:DB:580:U:O3'	38:DQ:30:VAL:HG13	2.14	0.47
23:DB:928:A:O2'	23:DB:929:U:H5'	2.15	0.47
25:DC:86:ARG:O	25:DC:88:ALA:N	2.47	0.47
26:DD:119:ALA:CB	26:DD:165:MET:HG2	2.44	0.47
27:DE:115:GLN:HG3	27:DE:116:ASP:N	2.27	0.47
27:DE:21:ARG:NH1	27:DE:23:PHE:HB3	2.24	0.47
52:DI:72:THR:HG23	52:DI:112:LYS:HD2	1.97	0.47
31:DJ:44:TYR:HD1	31:DJ:45:THR:H	1.51	0.47
36:DO:65:THR:HB	36:DO:66:GLY:H	1.56	0.47
36:DO:92:PHE:HE2	36:DO:94:ARG:HE	1.62	0.47
38:DQ:72:GLY:O	38:DQ:73:ILE:HG22	2.14	0.47
44:DX:49:ASP:HA	44:DX:52:ARG:HH12	1.79	0.47
1:AA:1246:A:H2'	1:AA:1247:U:O4'	2.15	0.47
1:AA:1348:U:H4'	8:AI:121:ARG:NH1	2.29	0.47
1:AA:202:G:H1'	1:AA:468:A:H8	1.79	0.47
1:AA:36:C:O2'	1:AA:37:U:H5'	2.15	0.47
1:AA:647:C:O2'	1:AA:648:A:H5'	2.14	0.47
1:AA:792:A:H1'	1:AA:794:A:N7	2.30	0.47
1:AA:923:A:H2'	1:AA:924:C:H6	1.79	0.47
1:AA:957:U:O2	1:AA:959:A:H8	1.98	0.47
2:AC:166:TRP:O	2:AC:167:TYR:HB2	2.14	0.47
3:AD:113:ALA:O	3:AD:117:VAL:HG23	2.15	0.47
8:AI:66:VAL:HG22	8:AI:67:LYS:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:963:G:H21	9:AJ:56:HIS:CE1	2.33	0.47
9:AJ:87:LEU:HB3	9:AJ:88:MET:CE	2.45	0.47
10:AK:21:HIS:CE1	10:AK:34:THR:HG21	2.50	0.47
10:AK:36:ARG:HG2	10:AK:37:GLN:N	2.30	0.47
13:AN:60:ARG:CD	13:AN:60:ARG:H	2.15	0.47
17:AR:64:LEU:HB2	17:AR:66:LEU:HG	1.97	0.47
19:AT:70:LYS:HZ3	19:AT:70:LYS:HB2	1.78	0.47
51:B4:22:VAL:HG11	51:B4:36:ARG:CZ	2.44	0.47
22:BA:113:C:H2'	22:BA:114:C:O4'	2.14	0.47
22:BA:30:C:H2'	22:BA:30:C:O2	2.15	0.47
23:BB:1707:G:H2'	23:BB:1708:C:H6	1.79	0.47
23:BB:1919:A:H2'	23:BB:1920:C:H5'	1.97	0.47
23:BB:2616:C:H2'	23:BB:2617:U:H6	1.79	0.47
23:BB:2630:G:H2'	23:BB:2631:G:C8	2.50	0.47
23:BB:2801:G:H2'	23:BB:2802:G:C8	2.49	0.47
23:BB:2862:G:H2'	23:BB:2863:C:H6	1.79	0.47
23:BB:32:C:H2'	23:BB:33:C:H5''	1.96	0.47
23:BB:459:U:H5'	49:B2:41:ARG:NE	2.27	0.47
23:BB:636:G:N7	33:BL:109:LYS:HE2	2.30	0.47
25:BC:38:LYS:HD3	25:BC:39:SER:O	2.14	0.47
26:BD:37:VAL:HG23	26:BD:42:ASN:ND2	2.30	0.47
23:BB:2786:U:H4'	26:BD:66:GLY:HA2	1.96	0.47
26:BD:83:ARG:HG2	26:BD:84:LEU:H	1.80	0.47
27:BE:143:LEU:HB3	27:BE:146:VAL:HG21	1.95	0.47
27:BE:148:ILE:HG22	27:BE:185:LYS:H	1.80	0.47
30:BH:72:ILE:O	30:BH:72:ILE:HG23	2.14	0.47
31:BJ:72:LYS:HG3	31:BJ:90:GLU:HG3	1.97	0.47
33:BL:82:LEU:CD2	33:BL:92:LEU:HD12	2.44	0.47
35:BN:32:GLU:HA	35:BN:115:LEU:HD21	1.95	0.47
38:BQ:95:ALA:C	38:BQ:97:ILE:N	2.67	0.47
42:BU:8:ASP:HB3	42:BU:25:LYS:HE2	1.95	0.47
43:BW:11:ASN:HD22	43:BW:13:ARG:NH2	2.13	0.47
45:BY:3:THR:HG21	45:BY:43:ILE:HB	1.95	0.47
1:CA:1144:G:N2	1:CA:1146:A:H62	2.12	0.47
1:CA:1157:A:H4'	1:CA:1158:C:O5'	2.14	0.47
1:CA:1381:U:O2'	6:CG:78:ARG:HB3	2.15	0.47
1:CA:1432:G:OP1	37:DP:105:LYS:HB2	2.15	0.47
3:CD:57:LYS:HD3	3:CD:58:GLN:N	2.29	0.47
9:CJ:55:PRO:O	9:CJ:56:HIS:HB3	2.14	0.47
13:CN:45:LEU:C	13:CN:47:LEU:H	2.18	0.47
50:D3:29:ARG:O	50:D3:31:ILE:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:D3:29:ARG:O	50:D3:30:HIS:C	2.53	0.47
23:DB:1241:A:O4'	23:DB:1241:A:N3	2.48	0.47
23:DB:1511:G:H2'	23:DB:1512:C:C6	2.50	0.47
23:DB:1544:A:H2'	23:DB:1545:A:C8	2.49	0.47
23:DB:1683:U:H2'	23:DB:1684:G:H8	1.80	0.47
23:DB:1824:G:H2'	23:DB:1825:U:C6	2.50	0.47
23:DB:2247:A:O2'	23:DB:2248:C:H5'	2.15	0.47
23:DB:2309:A:C6	23:DB:2310:C:N4	2.83	0.47
23:DB:2386:A:C4'	43:DW:38:ARG:HB2	2.44	0.47
23:DB:2760:C:H2'	23:DB:2760:C:O2	2.13	0.47
23:DB:2773:C:O2'	23:DB:2774:C:H5'	2.15	0.47
23:DB:967:U:H2'	23:DB:968:C:H6	1.78	0.47
23:DB:8:C:O2'	23:DB:9:G:H5'	2.15	0.47
25:DC:163:ILE:C	25:DC:164:VAL:HG22	2.34	0.47
27:DE:46:GLN:NE2	27:DE:48:THR:HB	2.29	0.47
28:DF:34:THR:O	28:DF:35:LEU:HD23	2.15	0.47
28:DF:37:MET:HB2	28:DF:86:CYS:SG	2.53	0.47
31:DJ:11:VAL:O	31:DJ:11:VAL:HG22	2.13	0.47
31:DJ:4:PHE:CD1	31:DJ:5:THR:N	2.83	0.47
31:DJ:93:ILE:HG22	31:DJ:93:ILE:O	2.15	0.47
34:DM:44:ARG:HB2	34:DM:91:TYR:HE2	1.80	0.47
35:DN:56:LYS:HG3	35:DN:57:THR:N	2.27	0.47
38:DQ:82:LEU:HD12	38:DQ:112:ALA:HB2	1.95	0.47
39:DR:49:ILE:HG13	39:DR:51:VAL:CG2	2.42	0.47
41:DT:55:VAL:CG2	41:DT:86:THR:H	2.28	0.47
44:DX:47:ARG:HE	44:DX:47:ARG:HA	1.79	0.47
1:AA:1423:G:H2'	1:AA:1424:U:O4'	2.14	0.47
1:AA:470:C:O2'	1:AA:471:U:H5'	2.14	0.47
1:AA:563:A:H2'	1:AA:567:G:C8	2.50	0.47
1:AA:824:G:O2'	1:AA:825:A:H5'	2.14	0.47
1:AA:940:C:H2'	1:AA:941:G:H8	1.80	0.47
1:AA:979:C:H2'	1:AA:980:C:O4'	2.14	0.47
2:AC:100:ILE:O	2:AC:100:ILE:HG23	2.14	0.47
2:AC:134:LYS:O	2:AC:138:GLN:HB2	2.14	0.47
1:AA:620:C:N1	3:AD:131:ILE:HD13	2.29	0.47
11:AL:30:ARG:HB3	11:AL:57:THR:CG2	2.45	0.47
14:AO:84:LEU:HD22	14:AO:86:LEU:HD11	1.96	0.47
15:AP:6:LEU:HB3	15:AP:17:TYR:HB3	1.96	0.47
48:B1:33:LEU:O	48:B1:33:LEU:HD12	2.14	0.47
23:BB:1083:U:H1'	23:BB:1086:A:N6	2.30	0.47
23:BB:1351:C:H2'	23:BB:1352:U:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1511:G:H2'	23:BB:1512:C:C6	2.50	0.47
23:BB:1651:G:H4'	35:BN:39:PRO:HG2	1.97	0.47
23:BB:1720:U:O2'	23:BB:1721:G:H5'	2.15	0.47
23:BB:1706:C:C2	23:BB:1757:A:H5'	2.49	0.47
23:BB:2041:U:H2'	23:BB:2042:A:H8	1.73	0.47
23:BB:2051:A:H5'	26:BD:145:SER:HB3	1.97	0.47
23:BB:2052:A:N9	26:BD:155:VAL:HG23	2.29	0.47
23:BB:2192:U:H2'	23:BB:2193:G:O4'	2.15	0.47
23:BB:2756:U:H1'	23:BB:2757:A:H5''	1.97	0.47
23:BB:2810:A:H2'	23:BB:2811:G:O4'	2.15	0.47
23:BB:282:A:H2'	23:BB:283:G:H8	1.76	0.47
23:BB:41:C:O2'	23:BB:42:A:H5'	2.13	0.47
23:BB:825:A:H2'	23:BB:826:U:O4'	2.14	0.47
25:BC:248:GLY:C	25:BC:250:GLN:H	2.18	0.47
25:BC:53:ILE:HG12	25:BC:218:THR:HA	1.96	0.47
27:BE:149:ILE:HG12	27:BE:186:VAL:HG22	1.97	0.47
27:BE:6:LYS:HE3	27:BE:118:LEU:HB2	1.96	0.47
28:BF:35:LEU:HD12	28:BF:35:LEU:N	2.30	0.47
29:BG:22:VAL:HG23	29:BG:33:THR:HG23	1.97	0.47
33:BL:77:ILE:HA	33:BL:101:ILE:CD1	2.45	0.47
37:BP:4:ILE:O	37:BP:7:LEU:HB3	2.15	0.47
37:BP:8:GLU:HA	37:BP:11:GLN:HG2	1.95	0.47
40:BS:33:LEU:HB3	40:BS:37:THR:HB	1.96	0.47
42:BU:13:LEU:CD2	42:BU:14:THR:H	2.28	0.47
24:BV:53:LYS:HD2	24:BV:54:ALA:N	2.29	0.47
1:CA:1009:U:H1'	1:CA:1021:A:C6	2.50	0.47
1:CA:1238:A:N3	1:CA:1241:G:H1'	2.30	0.47
1:CA:392:C:C2	1:CA:393:A:C8	3.03	0.47
1:CA:397:A:H5'	1:CA:398:U:OP1	2.13	0.47
1:CA:60:A:H1'	1:CA:61:G:O4'	2.14	0.47
1:CA:932:C:H2'	1:CA:933:G:C8	2.50	0.47
2:CC:66:THR:HG23	2:CC:101:ASN:HB2	1.97	0.47
2:CC:151:GLU:HG3	2:CC:166:TRP:HB2	1.97	0.47
5:CF:55:HIS:N	5:CF:55:HIS:ND1	2.62	0.47
6:CG:61:PHE:O	6:CG:62:GLU:C	2.51	0.47
12:CM:21:ILE:HB	12:CM:24:VAL:HG22	1.95	0.47
19:CT:78:LEU:O	19:CT:82:ILE:HG23	2.14	0.47
22:DA:50:A:OP1	36:DO:68:LYS:HB2	2.15	0.47
22:DA:82:U:H5''	45:DY:16:LEU:CD1	2.44	0.47
23:DB:105:C:H2'	23:DB:106:C:C6	2.50	0.47
23:DB:1656:C:H5''	26:DD:141:ARG:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1666:G:C2'	23:DB:1667:G:H5'	2.45	0.47
23:DB:2637:U:C2'	23:DB:2638:G:H5'	2.44	0.47
23:DB:2810:A:H2'	23:DB:2811:G:O4'	2.15	0.47
23:DB:356:G:O2'	23:DB:357:C:H5'	2.14	0.47
23:DB:852:U:H2'	23:DB:853:C:H6	1.79	0.47
25:DC:136:VAL:O	25:DC:138:SER:N	2.48	0.47
25:DC:42:ARG:NE	25:DC:42:ARG:HA	2.30	0.47
26:DD:122:VAL:HA	26:DD:128:ARG:HD2	1.96	0.47
26:DD:34:VAL:HG11	26:DD:50:VAL:HG23	1.97	0.47
26:DD:46:ARG:CA	26:DD:82:PHE:HA	2.44	0.47
26:DD:73:VAL:HB	26:DD:91:THR:CB	2.44	0.47
29:DG:37:ASN:N	29:DG:40:VAL:HG21	2.29	0.47
30:DH:94:ILE:HG23	30:DH:98:ASP:CG	2.35	0.47
52:DI:12:VAL:HG13	52:DI:41:PHE:CE2	2.50	0.47
33:DL:61:LEU:N	33:DL:62:PRO:CD	2.78	0.47
35:DN:12:ARG:N	35:DN:12:ARG:HD3	2.29	0.47
37:DP:4:ILE:O	37:DP:4:ILE:HG22	2.15	0.47
38:DQ:47:ARG:HH12	38:DQ:50:ARG:CG	2.27	0.47
39:DR:92:TRP:HE3	39:DR:93:PHE:N	2.13	0.47
43:DW:18:LYS:NZ	43:DW:18:LYS:HB2	2.30	0.47
43:DW:42:THR:HB	43:DW:75:ASN:HB3	1.94	0.47
44:DX:23:ARG:O	44:DX:26:PHE:HB2	2.14	0.47
44:DX:27:ASN:C	44:DX:29:ARG:H	2.18	0.47
44:DX:28:LEU:HD22	44:DX:42:LEU:HG	1.95	0.47
45:DY:1:ALA:H2	45:DY:43:ILE:HB	1.80	0.47
45:DY:6:ILE:HG12	45:DY:35:VAL:O	2.15	0.47
1:AA:1260:G:OP1	1:AA:1284:C:H4'	2.14	0.47
1:AA:1281:C:C5'	1:AA:1282:C:H5	2.27	0.47
1:AA:1300:G:O2'	1:AA:1301:U:P	2.73	0.47
1:AA:357:G:OP1	1:AA:367:U:H6	1.98	0.47
1:AA:392:C:C2	1:AA:393:A:C8	3.02	0.47
1:AA:408:A:H3'	1:AA:409:U:C6	2.47	0.47
1:AA:411:A:OP2	3:AD:25:ARG:NH2	2.47	0.47
1:AA:836:G:H2'	1:AA:837:U:H6	1.79	0.47
1:AA:893:C:H2'	1:AA:894:G:C8	2.49	0.47
1:AA:926:G:H3'	1:AA:1505:G:H21	1.80	0.47
20:AB:22:TRP:CG	20:AB:23:ASN:N	2.82	0.47
2:AC:38:VAL:O	2:AC:42:LEU:HD23	2.15	0.47
3:AD:25:ARG:CD	3:AD:26:ALA:H	2.24	0.47
4:AE:104:ILE:HG23	4:AE:111:ARG:HH12	1.80	0.47
8:AI:64:ILE:HG22	8:AI:65:THR:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AK:83:VAL:HG22	10:AK:106:ILE:HD11	1.97	0.47
11:AL:49:ARG:HH12	11:AL:88:ASP:CB	2.27	0.47
15:AP:73:ALA:HB1	15:AP:77:GLU:OE2	2.15	0.47
16:AQ:10:ARG:CZ	16:AQ:56:ASP:H	2.27	0.47
18:AS:51:HIS:HA	18:AS:55:GLN:O	2.15	0.47
23:BB:1010:A:N3	23:BB:1153:C:H1'	2.29	0.47
23:BB:1367:A:H2'	23:BB:1368:G:H5'	1.95	0.47
23:BB:1424:G:H2'	23:BB:1425:G:O4'	2.15	0.47
23:BB:1750:G:H2'	23:BB:1751:U:C6	2.50	0.47
23:BB:1775:U:H2'	23:BB:1776:G:O4'	2.14	0.47
23:BB:189:G:P	46:BZ:12:ILE:HG21	2.55	0.47
23:BB:2081:U:C5'	46:BZ:23:LYS:HD2	2.44	0.47
23:BB:213:A:H2'	23:BB:214:G:C8	2.50	0.47
23:BB:2800:A:N3	23:BB:2801:G:H1'	2.30	0.47
23:BB:309:A:H1'	23:BB:329:G:N3	2.30	0.47
23:BB:528:A:N1	23:BB:2043:C:C5'	2.78	0.47
23:BB:1789:A:OP1	25:BC:219:VAL:HG12	2.14	0.47
25:BC:3:VAL:HG12	25:BC:4:LYS:N	2.24	0.47
30:BH:6:LEU:HD21	30:BH:50:ARG:HD3	1.97	0.47
23:BB:1081:U:H5'	52:BI:126:ARG:HD2	1.96	0.47
31:BJ:12:LYS:HZ2	31:BJ:41:LYS:HG3	1.79	0.47
33:BL:119:PRO:O	33:BL:119:PRO:CD	2.62	0.47
39:BR:31:GLU:O	39:BR:32:THR:HB	2.15	0.47
39:BR:59:ILE:HG13	39:BR:59:ILE:O	2.14	0.47
39:BR:70:GLU:O	39:BR:71:LYS:HB2	2.14	0.47
41:BT:24:MET:HE3	41:BT:29:THR:H	1.79	0.47
1:CA:1085:U:H3'	1:CA:1086:U:C6	2.50	0.47
1:CA:1229:A:H2'	1:CA:1230:C:H6	1.79	0.47
1:CA:1300:G:C2'	1:CA:1301:U:OP2	2.62	0.47
1:CA:1301:U:O2'	1:CA:1302:C:OP1	2.25	0.47
1:CA:1512:U:H2'	1:CA:1513:A:C8	2.50	0.47
1:CA:284:C:H2'	1:CA:285:C:H6	1.79	0.47
1:CA:382:A:O2'	1:CA:383:A:H5'	2.15	0.47
1:CA:434:U:H2'	1:CA:434:U:O2	2.14	0.47
1:CA:476:U:O2'	1:CA:477:C:H5'	2.14	0.47
1:CA:719:C:O2'	17:CR:37:LYS:HB2	2.15	0.47
1:CA:890:G:O2'	1:CA:906:A:N6	2.48	0.47
2:CC:178:ARG:HG2	2:CC:205:GLU:O	2.15	0.47
3:CD:25:ARG:CB	3:CD:25:ARG:HH11	2.19	0.47
4:CE:113:VAL:HG21	4:CE:139:THR:HG21	1.96	0.47
7:CH:107:LYS:HB3	7:CH:107:LYS:NZ	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CI:56:MET:O	8:CI:58:GLU:HG2	2.14	0.47
9:CJ:7:ARG:HD2	9:CJ:101:SER:OG	2.14	0.47
11:CL:35:ARG:NE	11:CL:36:VAL:H	2.13	0.47
11:CL:81:ILE:HD13	11:CL:96:THR:HG22	1.96	0.47
17:CR:59:LYS:O	17:CR:63:TYR:HD2	1.97	0.47
21:CU:41:THR:O	21:CU:45:LYS:HB2	2.15	0.47
47:D0:29:VAL:HG22	47:D0:30:ASP:N	2.28	0.47
47:D0:36:LYS:HG2	47:D0:37:HIS:O	2.15	0.47
49:D2:18:PHE:CE2	49:D2:44:VAL:HB	2.50	0.47
50:D3:25:HIS:O	50:D3:27:ASN:N	2.48	0.47
23:DB:1060:U:OP1	52:DI:75:ALA:HB3	2.15	0.47
23:DB:1332:G:N3	23:DB:1332:G:H2'	2.29	0.47
23:DB:1351:C:H2'	23:DB:1352:U:C6	2.49	0.47
23:DB:143:C:H2'	23:DB:144:A:C8	2.50	0.47
23:DB:1541:C:H2'	23:DB:1542:U:C6	2.50	0.47
23:DB:2462:C:H2'	23:DB:2463:C:C6	2.49	0.47
23:DB:2597:G:OP1	25:DC:239:PHE:CG	2.68	0.47
23:DB:2710:C:H2'	23:DB:2711:A:H8	1.80	0.47
23:DB:2864:G:H2'	23:DB:2865:U:C6	2.49	0.47
25:DC:226:PRO:HG3	25:DC:232:GLY:C	2.35	0.47
25:DC:52:HIS:O	25:DC:53:ILE:HD12	2.15	0.47
26:DD:112:THR:HB	26:DD:167:ASN:O	2.14	0.47
27:DE:142:ALA:C	27:DE:143:LEU:HD13	2.34	0.47
27:DE:147:LEU:HB3	27:DE:167:VAL:HG22	1.96	0.47
28:DF:39:VAL:HG13	28:DF:84:ILE:CB	2.45	0.47
28:DF:32:LYS:CB	28:DF:91:ARG:HB3	2.33	0.47
29:DG:23:ILE:O	29:DG:33:THR:HA	2.14	0.47
52:DI:45:THR:O	52:DI:48:ILE:HG22	2.14	0.47
52:DI:69:VAL:HG23	52:DI:69:VAL:O	2.15	0.47
33:DL:142:ILE:HG12	33:DL:144:GLU:OE1	2.15	0.47
33:DL:59:ARG:O	33:DL:60:ARG:HD2	2.14	0.47
41:DT:21:SER:O	41:DT:24:MET:HB2	2.15	0.47
41:DT:62:VAL:O	41:DT:63:VAL:HB	2.15	0.47
42:DU:15:GLY:C	42:DU:17:ASP:H	2.18	0.47
46:DZ:1:MET:HA	46:DZ:9:TYR:CG	2.50	0.47
46:DZ:59:ARG:O	46:DZ:61:ASN:N	2.41	0.47
1:AA:1117:A:H4'	8:AI:105:ARG:NH1	2.30	0.47
1:AA:513:C:H2'	1:AA:514:C:H6	1.79	0.47
1:AA:768:A:H4'	1:AA:1523:G:N2	2.30	0.47
1:AA:803:G:H2'	1:AA:804:U:C6	2.50	0.47
20:AB:15:PHE:HD1	20:AB:16:GLY:H	1.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1112:C:O2	2:AC:178:ARG:N	2.48	0.47
2:AC:76:ILE:O	2:AC:83:VAL:HG23	2.14	0.47
2:AC:90:VAL:HA	2:AC:93:ILE:CG2	2.44	0.47
3:AD:54:LEU:O	3:AD:57:LYS:HB3	2.14	0.47
3:AD:71:PHE:HA	3:AD:74:TYR:HD2	1.80	0.47
12:AM:16:ILE:HG23	12:AM:17:ALA:N	2.30	0.47
1:AA:1226:C:P	12:AM:89:ARG:HH12	2.37	0.47
19:AT:57:VAL:HG12	19:AT:57:VAL:O	2.15	0.47
1:AA:263:A:P	19:AT:73:ARG:HH22	2.38	0.47
50:B3:24:LYS:NZ	50:B3:29:ARG:HH12	2.13	0.47
23:BB:125:A:H4'	49:B2:13:ASN:CG	2.34	0.47
23:BB:1710:G:O2'	23:BB:1711:A:H5'	2.14	0.47
23:BB:2023:C:H4'	23:BB:2617:U:O3'	2.15	0.47
23:BB:2038:G:H2'	23:BB:2039:U:C6	2.49	0.47
23:BB:236:C:O2'	23:BB:237:C:H5'	2.14	0.47
23:BB:2519:U:H5'	23:BB:2567:G:N2	2.30	0.47
23:BB:2603:G:O2'	23:BB:2604:U:H5'	2.15	0.47
23:BB:2773:C:O2'	23:BB:2774:C:H5'	2.15	0.47
23:BB:527:C:C4	23:BB:2779:U:H2'	2.50	0.47
23:BB:55:G:H2'	23:BB:56:A:C8	2.50	0.47
23:BB:599:A:O2'	23:BB:600:G:H5'	2.15	0.47
23:BB:664:G:O2'	23:BB:665:U:H5'	2.15	0.47
25:BC:224:MET:HB3	25:BC:233:GLY:H	1.80	0.47
29:BG:126:THR:C	29:BG:128:THR:H	2.17	0.47
29:BG:32:LEU:HD21	29:BG:74:MET:HB3	1.97	0.47
30:BH:121:VAL:HG11	30:BH:128:HIS:CD2	2.49	0.47
52:BI:83:ALA:N	52:BI:100:ILE:HD11	2.29	0.47
32:BK:73:ASP:OD1	32:BK:74:GLY:N	2.48	0.47
33:BL:39:LYS:HB2	33:BL:46:VAL:HG13	1.96	0.47
34:BM:103:TYR:O	34:BM:105:MET:N	2.48	0.47
34:BM:116:ALA:O	34:BM:119:LEU:HD23	2.15	0.47
34:BM:3:GLN:C	34:BM:5:LYS:N	2.66	0.47
36:BO:26:LEU:HG	36:BO:39:VAL:HG22	1.96	0.47
44:BX:22:LEU:CD2	44:BX:25:GLN:HB3	2.45	0.47
1:CA:1225:A:H5'	12:CM:101:THR:HG1	1.80	0.47
1:CA:197:A:H4'	1:CA:198:G:O5'	2.15	0.47
1:CA:252:U:H2'	1:CA:253:A:C8	2.49	0.47
1:CA:647:C:O2'	1:CA:648:A:H5'	2.15	0.47
1:CA:695:A:H5'	10:CK:52:ARG:NH2	2.28	0.47
20:CB:199:ILE:O	20:CB:199:ILE:HG13	2.15	0.47
12:CM:95:PRO:CD	12:CM:108:ARG:HG2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:CO:23:SER:HB3	14:CO:26:VAL:CG2	2.45	0.47
15:CP:53:ASP:O	15:CP:57:ILE:HG22	2.14	0.47
17:CR:20:ILE:O	17:CR:20:ILE:HG12	2.15	0.47
18:CS:14:LEU:HD12	18:CS:15:LEU:N	2.30	0.47
23:DB:129:C:H4'	23:DB:1348:C:O2'	2.15	0.47
23:DB:1355:G:O2'	23:DB:1356:G:H5'	2.14	0.47
23:DB:1427:A:H4'	23:DB:1428:C:O4'	2.14	0.47
23:DB:1710:G:O2'	23:DB:1711:A:H5'	2.14	0.47
23:DB:1720:U:O2'	23:DB:1721:G:H5'	2.15	0.47
23:DB:1882:U:O2'	23:DB:1883:U:H5'	2.14	0.47
23:DB:2093:G:O2'	23:DB:2094:A:H5'	2.14	0.47
23:DB:2722:G:H2'	23:DB:2723:C:C6	2.49	0.47
23:DB:2733:A:H3'	23:DB:2733:A:C8	2.50	0.47
23:DB:2801:G:H2'	23:DB:2802:G:C8	2.50	0.47
23:DB:543:G:C2'	23:DB:545:U:OP1	2.63	0.47
23:DB:5:A:H2'	23:DB:6:A:C8	2.50	0.47
23:DB:599:A:O2'	23:DB:600:G:H5'	2.15	0.47
23:DB:962:G:O2'	23:DB:963:U:H5'	2.15	0.47
25:DC:52:HIS:O	25:DC:53:ILE:CB	2.61	0.47
26:DD:50:VAL:HG13	26:DD:77:ARG:O	2.14	0.47
27:DE:123:LYS:NZ	27:DE:158:PHE:HA	2.30	0.47
27:DE:169:VAL:HG22	27:DE:171:ASP:H	1.80	0.47
28:DF:177:ARG:O	28:DF:178:LYS:HB2	2.15	0.47
28:DF:56:LEU:HD13	28:DF:88:VAL:HG21	1.97	0.47
29:DG:10:VAL:HG11	29:DG:44:HIS:CE1	2.50	0.47
29:DG:8:VAL:HG12	29:DG:49:LEU:HD12	1.96	0.47
29:DG:25:ILE:CD1	29:DG:75:VAL:HG22	2.45	0.47
30:DH:88:GLY:O	30:DH:124:THR:HA	2.15	0.47
31:DJ:120:ARG:C	31:DJ:121:LYS:HE3	2.35	0.47
23:DB:1665:A:H5''	32:DK:66:LYS:HB3	1.97	0.47
34:DM:62:LYS:H	34:DM:104:GLU:CB	2.22	0.47
35:DN:103:ARG:O	35:DN:104:ALA:HB3	2.15	0.47
36:DO:35:ILE:CG2	36:DO:74:VAL:HG21	2.44	0.47
37:DP:65:ASN:H	37:DP:71:ARG:CA	2.27	0.47
38:DQ:116:LEU:H	38:DQ:116:LEU:HD22	1.79	0.47
23:DB:1009:A:H5'	38:DQ:58:GLN:HE21	1.80	0.47
38:DQ:82:LEU:HG	38:DQ:88:GLU:CG	2.45	0.47
41:DT:40:LYS:HE3	41:DT:59:ASN:O	2.14	0.47
24:DV:65:VAL:O	24:DV:66:ASP:HB3	2.15	0.47
1:AA:1111:A:C2	2:AC:176:THR:HG23	2.50	0.47
1:AA:1212:U:H5'	1:AA:1213:A:O4'	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1406:U:H3'	1:AA:1407:C:H6	1.79	0.47
1:AA:284:C:H2'	1:AA:285:C:H6	1.79	0.47
1:AA:413:G:H5'	1:AA:414:A:OP1	2.14	0.47
1:AA:589:U:O2'	1:AA:590:U:H5'	2.14	0.47
1:AA:652:U:H1'	1:AA:653:U:C5	2.50	0.47
20:AB:9:LEU:HD22	20:AB:10:LYS:N	2.30	0.47
2:AC:115:VAL:CG1	2:AC:137:VAL:HG13	2.44	0.47
2:AC:53:ARG:HG2	2:AC:54:ILE:N	2.26	0.47
4:AE:105:ILE:HB	4:AE:123:LEU:CA	2.44	0.47
5:AF:51:ILE:O	5:AF:51:ILE:HG23	2.15	0.47
8:AI:66:VAL:HG21	8:AI:74:GLN:HG3	1.97	0.47
11:AL:35:ARG:HE	11:AL:35:ARG:CA	2.23	0.47
11:AL:58:ASN:H	11:AL:58:ASN:ND2	2.12	0.47
15:AP:67:ILE:CD1	15:AP:72:ALA:HA	2.45	0.47
16:AQ:79:GLU:HG3	16:AQ:80:LYS:N	2.31	0.47
18:AS:69:LYS:O	18:AS:72:GLU:HB2	2.15	0.47
19:AT:34:VAL:HG11	19:AT:78:LEU:HD13	1.97	0.47
1:AA:108:G:C6	19:AT:9:ARG:HG2	2.50	0.47
47:B0:42:ILE:HD11	47:B0:44:ALA:HB3	1.97	0.47
48:B1:31:GLU:O	48:B1:33:LEU:HG	2.15	0.47
50:B3:56:LEU:O	50:B3:57:VAL:HB	2.15	0.47
22:BA:105:G:O2'	22:BA:106:G:H5'	2.15	0.47
23:BB:1515:A:H5'	23:BB:1557:C:C5'	2.44	0.47
23:BB:173:A:H2'	23:BB:174:U:H6	1.80	0.47
23:BB:2299:U:H2'	23:BB:2300:C:C6	2.50	0.47
23:BB:2543:G:H2'	23:BB:2544:G:C8	2.50	0.47
23:BB:2822:G:N7	35:BN:1:MET:N	2.58	0.47
23:BB:327:G:H2'	23:BB:328:U:C6	2.50	0.47
23:BB:535:G:H3'	55:BB:3420:HOH:O	2.15	0.47
23:BB:657:U:H2'	23:BB:658:U:C6	2.50	0.47
23:BB:764:A:OP2	25:BC:213:ARG:NH2	2.47	0.47
26:BD:122:VAL:CB	26:BD:141:ARG:HH12	2.28	0.47
26:BD:29:VAL:HG12	26:BD:185:ASN:ND2	2.30	0.47
26:BD:45:TYR:CE2	26:BD:83:ARG:HG3	2.50	0.47
30:BH:30:LEU:HD13	30:BH:35:LYS:HD3	1.97	0.47
52:BI:29:GLN:HA	52:BI:29:GLN:HE21	1.80	0.47
31:BJ:63:ALA:HA	31:BJ:69:ARG:HD3	1.97	0.47
32:BK:99:ILE:N	32:BK:99:ILE:HD12	2.30	0.47
23:BB:661:A:O2'	33:BL:23:ILE:HG23	2.14	0.47
34:BM:21:ALA:O	34:BM:96:ILE:HG13	2.15	0.47
35:BN:103:ARG:HB2	35:BN:110:MET:SD	2.55	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:49:GLU:O	35:BN:52:ILE:HB	2.15	0.47
36:BO:84:GLU:OE1	36:BO:85:LYS:HG3	2.14	0.47
38:BQ:20:ALA:O	38:BQ:22:GLY:N	2.48	0.47
39:BR:5:PHE:HB3	39:BR:12:HIS:CG	2.50	0.47
41:BT:48:GLN:HG3	41:BT:49:LYS:N	2.30	0.47
44:BX:36:GLN:O	44:BX:37:LEU:HD23	2.15	0.47
44:BX:38:GLN:HG2	44:BX:39:GLN:N	2.29	0.47
46:BZ:49:ARG:O	46:BZ:51:VAL:HG23	2.14	0.47
1:CA:840:C:N3	1:CA:842:U:H4'	2.30	0.47
1:CA:992:U:H1'	1:CA:993:G:C2	2.49	0.47
2:CC:127:VAL:HB	2:CC:128:MET:CE	2.45	0.47
3:CD:13:ARG:HG3	3:CD:13:ARG:O	2.15	0.47
4:CE:45:VAL:O	4:CE:71:ILE:HG22	2.14	0.47
6:CG:100:MET:O	6:CG:104:VAL:HG23	2.14	0.47
6:CG:29:LEU:HD12	6:CG:42:VAL:HG22	1.96	0.47
1:CA:948:C:OP1	12:CM:105:ALA:HA	2.14	0.47
23:DB:1098:A:O5'	52:DI:3:LYS:CG	2.63	0.47
23:DB:1199:U:C2	23:DB:1200:C:C5	3.03	0.47
23:DB:1409:U:O2'	23:DB:1410:G:H5'	2.15	0.47
23:DB:2071:A:H2'	23:DB:2072:C:H6	1.76	0.47
23:DB:2300:C:H2'	23:DB:2301:C:H6	1.79	0.47
23:DB:299:A:H2	23:DB:319:G:N3	2.13	0.47
23:DB:346:A:N7	23:DB:347:A:H1'	2.30	0.47
23:DB:657:U:H2'	23:DB:658:U:C6	2.50	0.47
23:DB:802:A:H4'	55:DB:3289:HOH:O	2.15	0.47
23:DB:902:C:O2'	23:DB:903:C:H5'	2.15	0.47
23:DB:962:G:O2'	23:DB:2250:G:N2	2.48	0.47
25:DC:12:ARG:HH11	25:DC:18:VAL:CB	2.28	0.47
27:DE:88:ARG:HB3	27:DE:89:PRO:HD2	1.97	0.47
30:DH:26:ALA:C	30:DH:28:ASN:H	2.16	0.47
30:DH:68:ARG:HD3	30:DH:71:LYS:HD3	1.97	0.47
32:DK:20:MET:C	32:DK:41:ILE:HD12	2.34	0.47
34:DM:33:LEU:CD2	34:DM:124:LEU:HB2	2.44	0.47
35:DN:96:ARG:HG2	35:DN:96:ARG:HH11	1.79	0.47
37:DP:59:THR:HG1	37:DP:76:HIS:CG	2.33	0.47
40:DS:7:HIS:NE2	40:DS:10:ALA:HA	2.29	0.47
41:DT:68:LYS:HB2	41:DT:68:LYS:HZ2	1.77	0.47
42:DU:72:PHE:CD2	42:DU:74:ALA:HB3	2.50	0.47
1:AA:1432:G:C5'	37:BP:106:ALA:HB2	2.44	0.46
1:AA:1456:A:H2'	1:AA:1457:G:O4'	2.16	0.46
1:AA:216:U:H2'	1:AA:217:C:C6	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:455:G:H2'	1:AA:456:A:C8	2.51	0.46
1:AA:601:G:H2'	1:AA:602:A:H8	1.80	0.46
1:AA:691:G:H1'	1:AA:696:A:N6	2.30	0.46
1:AA:909:A:H2'	1:AA:910:C:O4'	2.14	0.46
3:AD:196:GLU:HA	3:AD:199:ILE:HD12	1.97	0.46
3:AD:24:VAL:HG23	3:AD:25:ARG:HD2	1.96	0.46
7:AH:64:TYR:CA	7:AH:70:VAL:HG23	2.44	0.46
7:AH:69:ALA:HB3	7:AH:72:GLU:OE2	2.15	0.46
8:AI:18:VAL:HG13	8:AI:64:ILE:HG12	1.96	0.46
9:AJ:73:LEU:O	9:AJ:74:VAL:HB	2.15	0.46
11:AL:1:ALA:HB1	11:AL:5:GLN:HB2	1.98	0.46
11:AL:89:LEU:HD22	11:AL:89:LEU:N	2.30	0.46
12:AM:70:ARG:NE	28:BF:112:ASP:OD2	2.49	0.46
13:AN:14:ALA:O	13:AN:18:LYS:HG2	2.15	0.46
13:AN:72:PHE:CE2	13:AN:77:GLY:HA2	2.50	0.46
14:AO:17:ASP:HB2	14:AO:18:ALA:H	1.57	0.46
16:AQ:10:ARG:CA	16:AQ:10:ARG:NE	2.77	0.46
50:B3:34:LYS:HB3	50:B3:35:LYS:NZ	2.30	0.46
23:BB:1029:A:H2'	23:BB:1030:C:O4'	2.15	0.46
23:BB:1156:A:N6	38:BQ:47:ARG:HD2	2.30	0.46
23:BB:1257:C:H5'	27:BE:78:TRP:CH2	2.49	0.46
23:BB:135:U:H2'	23:BB:136:G:N9	2.29	0.46
23:BB:1409:U:O2'	23:BB:1410:G:H5'	2.15	0.46
23:BB:528:A:C6	23:BB:2042:A:H2'	2.46	0.46
23:BB:2378:A:C2'	23:BB:2379:G:H5'	2.43	0.46
23:BB:241:A:OP1	23:BB:241:A:H8	1.98	0.46
23:BB:2769:U:O2'	23:BB:2770:G:H5'	2.16	0.46
23:BB:778:G:H5''	25:BC:48:ILE:HD12	1.97	0.46
23:BB:6:A:O2'	23:BB:7:G:H5'	2.15	0.46
23:BB:946:C:H2'	23:BB:947:A:H8	1.80	0.46
25:BC:160:TYR:HA	25:BC:193:GLU:HG2	1.96	0.46
27:BE:149:ILE:HD11	27:BE:186:VAL:HG13	1.97	0.46
27:BE:173:THR:O	27:BE:175:ILE:HG22	2.14	0.46
34:BM:108:VAL:O	34:BM:112:LEU:HD13	2.15	0.46
35:BN:27:SER:OG	35:BN:34:ILE:HG21	2.15	0.46
36:BO:66:GLY:O	36:BO:67:ASN:ND2	2.48	0.46
38:BQ:92:LYS:HD3	38:BQ:92:LYS:C	2.35	0.46
40:BS:24:ILE:HD13	40:BS:36:LEU:HD11	1.97	0.46
24:BV:82:TYR:CB	34:BM:36:VAL:HG12	2.45	0.46
43:BW:58:LEU:HA	43:BW:79:ILE:CG2	2.45	0.46
46:BZ:8:LYS:CD	46:BZ:9:TYR:N	2.70	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1028:C:H2'	1:CA:1029:U:C6	2.49	0.46
1:CA:1226:C:N4	12:CM:102:LYS:HZ3	2.12	0.46
1:CA:1363:A:H2'	1:CA:1363:A:N3	2.30	0.46
1:CA:216:U:H2'	1:CA:217:C:C6	2.50	0.46
1:CA:279:A:H5'	1:CA:281:G:O4'	2.15	0.46
1:CA:472:U:N3	1:CA:473:U:C4	2.82	0.46
1:CA:601:G:H2'	1:CA:602:A:C8	2.50	0.46
1:CA:831:A:O2'	1:CA:832:G:H5'	2.15	0.46
20:CB:18:GLN:HB2	20:CB:188:THR:OG1	2.14	0.46
2:CC:13:ILE:HG13	2:CC:177:LEU:HB3	1.98	0.46
7:CH:46:GLU:N	7:CH:63:LYS:HE3	2.30	0.46
8:CI:115:VAL:HG21	9:CJ:62:ARG:HG3	1.97	0.46
10:CK:127:ARG:HH11	10:CK:127:ARG:HG3	1.79	0.46
12:CM:105:ALA:O	12:CM:109:LYS:HG3	2.15	0.46
13:CN:10:VAL:HG12	13:CN:11:LYS:NZ	2.30	0.46
9:CJ:65:TYR:HB3	13:CN:95:LEU:HD11	1.96	0.46
47:D0:32:THR:HG21	47:D0:41:HIS:CE1	2.49	0.46
48:D1:7:LYS:HA	48:D1:26:LYS:HA	1.97	0.46
48:D1:15:GLY:HA3	48:D1:47:ILE:CG2	2.44	0.46
23:DB:1048:A:C5	23:DB:1111:A:C2	3.03	0.46
23:DB:1669:A:H2'	23:DB:1669:A:N3	2.31	0.46
23:DB:1936:A:H2	23:DB:1943:U:C5	2.33	0.46
23:DB:1997:C:H5'	26:DD:128:ARG:NH2	2.29	0.46
23:DB:1999:C:O2'	23:DB:2000:C:H5'	2.15	0.46
23:DB:2300:C:H2'	23:DB:2301:C:C6	2.51	0.46
23:DB:2420:C:H3'	50:D3:31:ILE:HG22	1.96	0.46
23:DB:356:G:H2'	23:DB:357:C:H6	1.80	0.46
23:DB:358:U:H2'	23:DB:359:G:H8	1.77	0.46
23:DB:533:G:N3	38:DQ:40:LYS:HG2	2.30	0.46
23:DB:957:C:O2'	23:DB:958:U:H5''	2.14	0.46
25:DC:38:LYS:O	25:DC:60:ALA:HA	2.14	0.46
25:DC:77:VAL:HG23	25:DC:113:ASP:HB2	1.97	0.46
27:DE:153:LEU:HD23	27:DE:172:ALA:N	2.31	0.46
28:DF:79:ARG:HB2	28:DF:82:TYR:CZ	2.50	0.46
52:DI:37:PHE:CE1	52:DI:58:ILE:HD11	2.50	0.46
31:DJ:100:VAL:HG13	31:DJ:101:ILE:HG12	1.96	0.46
33:DL:108:ALA:HB3	33:DL:125:LEU:HB3	1.97	0.46
33:DL:123:ARG:HA	33:DL:142:ILE:CA	2.45	0.46
34:DM:4:PRO:HB2	34:DM:69:PRO:CG	2.45	0.46
36:DO:86:GLY:C	36:DO:88:LYS:H	2.17	0.46
31:DJ:43:GLU:CA	38:DQ:63:ARG:HH12	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DS:4:ILE:HG12	40:DS:106:VAL:HG12	1.97	0.46
40:DS:7:HIS:NE2	40:DS:46:LEU:HD22	2.30	0.46
23:DB:64:A:H5'	41:DT:76:ARG:HD2	1.95	0.46
42:DU:90:LYS:O	42:DU:91:LYS:HB2	2.16	0.46
24:DV:69:GLU:O	24:DV:70:ILE:HG23	2.15	0.46
1:AA:1390:U:H2'	1:AA:1391:U:H6	1.79	0.46
1:AA:239:U:H5''	1:AA:239:U:H6	1.80	0.46
1:AA:478:A:H2'	1:AA:479:U:O4'	2.15	0.46
1:AA:393:A:H5'	1:AA:483:C:O2'	2.15	0.46
1:AA:691:G:H2'	1:AA:692:U:C6	2.50	0.46
1:AA:754:C:H3'	1:AA:754:C:O2	2.15	0.46
1:AA:889:A:H61	1:AA:907:A:H3'	1.80	0.46
20:AB:115:ASP:O	20:AB:119:GLN:HG2	2.15	0.46
9:AJ:88:MET:HB2	9:AJ:89:ARG:HH22	1.81	0.46
1:AA:676:A:H1'	10:AK:116:PRO:HB3	1.96	0.46
10:AK:126:ARG:CB	21:AU:33:ARG:HD2	2.39	0.46
49:B2:14:ARG:HD3	49:B2:14:ARG:O	2.14	0.46
23:BB:1190:G:OP1	33:BL:41:ARG:NE	2.48	0.46
23:BB:1526:C:H2'	23:BB:1527:G:O4'	2.15	0.46
23:BB:2263:C:H2'	23:BB:2264:C:C6	2.50	0.46
23:BB:2477:U:H4'	23:BB:2479:U:O4	2.16	0.46
23:BB:2733:A:H3'	23:BB:2733:A:C8	2.50	0.46
23:BB:834:G:H2'	23:BB:835:C:H6	1.79	0.46
25:BC:137:GLY:C	25:BC:139:THR:H	2.18	0.46
25:BC:32:LEU:HD13	25:BC:61:TYR:CE1	2.50	0.46
25:BC:53:ILE:HG21	25:BC:218:THR:HG23	1.95	0.46
25:BC:59:GLN:HB3	25:BC:60:ALA:H	1.51	0.46
27:BE:27:LEU:HD13	27:BE:27:LEU:C	2.36	0.46
23:BB:443:A:H5''	27:BE:44:ARG:HH12	1.80	0.46
28:BF:115:GLY:HA3	28:BF:177:ARG:N	2.31	0.46
30:BH:73:ASN:C	30:BH:75:LEU:H	2.19	0.46
52:BI:21:PRO:CB	52:BI:22:PRO:HD3	2.42	0.46
52:BI:44:LYS:O	52:BI:48:ILE:HG13	2.15	0.46
31:BJ:102:GLU:OE2	31:BJ:124:VAL:HG11	2.16	0.46
32:BK:114:LYS:O	32:BK:117:SER:HB3	2.15	0.46
32:BK:98:ARG:HA	32:BK:118:LEU:CD2	2.45	0.46
35:BN:45:ARG:N	35:BN:45:ARG:HD2	2.30	0.46
36:BO:6:ALA:HB1	36:BO:10:ARG:NH1	2.30	0.46
36:BO:57:ALA:HB3	36:BO:62:LEU:HA	1.97	0.46
37:BP:29:VAL:O	37:BP:30:TRP:C	2.53	0.46
38:BQ:105:PHE:O	38:BQ:109:VAL:HG13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BQ:85:ALA:HB3	38:BQ:88:GLU:O	2.15	0.46
39:BR:90:ARG:HH12	39:BR:91:GLN:C	2.18	0.46
40:BS:89:ALA:O	40:BS:90:LYS:HD2	2.16	0.46
41:BT:44:LYS:HB2	41:BT:44:LYS:NZ	2.30	0.46
24:BV:66:ASP:C	24:BV:68:LYS:H	2.18	0.46
43:BW:23:LYS:HG3	43:BW:24:ARG:N	2.30	0.46
1:CA:1049:U:H1'	1:CA:1201:A:N7	2.30	0.46
1:CA:1118:U:H2'	1:CA:1119:C:C6	2.47	0.46
1:CA:1368:A:O2'	1:CA:1369:C:H5'	2.15	0.46
1:CA:547:A:H4'	1:CA:548:G:O5'	2.16	0.46
1:CA:782:A:H2'	1:CA:783:C:O4'	2.14	0.46
2:CC:190:THR:HG22	2:CC:191:THR:N	2.31	0.46
3:CD:59:LYS:O	3:CD:63:ILE:HG22	2.15	0.46
16:CQ:44:HIS:HB2	16:CQ:69:THR:O	2.15	0.46
16:CQ:74:LEU:HD13	16:CQ:74:LEU:C	2.36	0.46
19:CT:67:HIS:CG	19:CT:68:LYS:H	2.31	0.46
48:D1:13:SER:HB3	48:D1:49:LYS:HZ3	1.79	0.46
22:DA:13:G:C2'	22:DA:14:U:H5''	2.45	0.46
23:DB:106:C:H2'	23:DB:107:G:C8	2.51	0.46
23:DB:146:A:H2'	23:DB:147:C:H6	1.80	0.46
23:DB:1720:U:C2'	23:DB:1721:G:H5'	2.46	0.46
23:DB:1735:A:H2'	23:DB:1736:U:C6	2.50	0.46
23:DB:1946:U:C2	23:DB:1947:C:C5	3.03	0.46
23:DB:2228:G:H2'	23:DB:2229:U:H6	1.77	0.46
23:DB:2502:G:H5'	23:DB:2503:A:C5'	2.36	0.46
23:DB:2515:C:H2'	23:DB:2516:A:H8	1.81	0.46
23:DB:2700:A:H2'	23:DB:2701:U:C6	2.51	0.46
23:DB:286:U:H2'	23:DB:287:G:C8	2.51	0.46
23:DB:2896:C:H2'	23:DB:2897:U:H6	1.80	0.46
23:DB:448:U:O4'	23:DB:448:U:OP2	2.34	0.46
23:DB:753:A:H2'	23:DB:754:U:H6	1.78	0.46
23:DB:931:U:N3	23:DB:1166:G:N2	2.63	0.46
23:DB:1820:U:O2'	25:DC:157:ALA:HB3	2.15	0.46
30:DH:83:LYS:O	30:DH:90:LEU:HA	2.14	0.46
33:DL:89:VAL:O	33:DL:89:VAL:HG13	2.16	0.46
35:DN:8:ARG:HD2	35:DN:46:ARG:CZ	2.44	0.46
36:DO:21:LEU:CD2	36:DO:22:GLY:H	2.23	0.46
37:DP:38:ARG:CG	37:DP:39:LEU:N	2.75	0.46
37:DP:54:LEU:N	37:DP:58:PHE:HD1	2.13	0.46
39:DR:58:VAL:O	39:DR:58:VAL:HG13	2.14	0.46
39:DR:76:LYS:O	39:DR:77:PHE:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DS:41:LYS:HG2	40:DS:43:ALA:HB3	1.97	0.46
43:DW:33:GLY:O	43:DW:66:VAL:HG23	2.15	0.46
44:DX:4:LYS:HG3	44:DX:7:ARG:HE	1.81	0.46
45:DY:2:LYS:N	45:DY:37:ARG:HB2	2.30	0.46
23:DB:2090:A:H2'	46:DZ:49:ARG:CZ	2.45	0.46
1:AA:1005:A:H3'	1:AA:1006:G:C8	2.48	0.46
1:AA:1050:G:C2'	1:AA:1051:C:H5'	2.44	0.46
1:AA:1184:G:C2	1:AA:1185:G:C8	3.03	0.46
1:AA:1234:C:H2'	1:AA:1235:U:C6	2.50	0.46
1:AA:1243:C:H2'	1:AA:1244:G:C8	2.51	0.46
1:AA:252:U:H2'	1:AA:253:A:C8	2.50	0.46
1:AA:279:A:H5'	1:AA:281:G:O4'	2.15	0.46
1:AA:547:A:H4'	1:AA:548:G:O5'	2.16	0.46
1:AA:598:U:H2'	1:AA:599:C:C6	2.51	0.46
1:AA:611:C:H2'	1:AA:612:C:H6	1.80	0.46
1:AA:924:C:H2'	1:AA:925:G:H8	1.81	0.46
20:AB:61:SER:HA	20:AB:224:ARG:CA	2.43	0.46
20:AB:9:LEU:H	20:AB:9:LEU:CD1	2.27	0.46
6:AG:91:ARG:HD3	6:AG:91:ARG:H	1.80	0.46
8:AI:18:VAL:HG21	8:AI:82:ILE:CG1	2.45	0.46
12:AM:3:ILE:HA	12:AM:56:ARG:CG	2.41	0.46
16:AQ:26:ARG:HE	16:AQ:39:ARG:NH1	2.14	0.46
16:AQ:65:PRO:HA	16:AQ:71:SER:OG	2.15	0.46
19:AT:60:GLN:NE2	19:AT:61:ALA:H	2.12	0.46
19:AT:66:ILE:HG23	19:AT:70:LYS:CD	2.45	0.46
47:B0:35:GLU:HG2	47:B0:36:LYS:N	2.30	0.46
48:B1:49:LYS:N	48:B1:49:LYS:HD2	2.30	0.46
23:BB:1179:G:O2'	23:BB:1180:U:H5'	2.16	0.46
23:BB:1186:G:H2'	23:BB:1187:G:O4'	2.16	0.46
23:BB:1541:C:H2'	23:BB:1542:U:C6	2.50	0.46
23:BB:1666:G:C2'	23:BB:1667:G:H5'	2.45	0.46
23:BB:1824:G:H2'	23:BB:1825:U:H6	1.81	0.46
23:BB:1947:C:O2'	23:BB:1948:G:H5'	2.15	0.46
23:BB:2199:A:H3'	23:BB:2200:C:H6	1.80	0.46
23:BB:2317:A:N6	23:BB:2318:G:C4	2.83	0.46
23:BB:2320:U:H3'	23:BB:2321:U:C5'	2.46	0.46
23:BB:2714:G:H2'	23:BB:2715:C:H6	1.80	0.46
23:BB:372:G:O2'	23:BB:373:U:P	2.73	0.46
23:BB:633:A:H2'	23:BB:634:C:O4'	2.15	0.46
25:BC:128:THR:HA	25:BC:190:THR:CA	2.40	0.46
25:BC:155:ARG:O	25:BC:156:SER:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:161:SER:HB3	28:BF:164:GLU:HB2	1.97	0.46
29:BG:17:LYS:O	29:BG:23:ILE:HD12	2.14	0.46
31:BJ:98:GLU:O	31:BJ:100:VAL:N	2.48	0.46
33:BL:78:ARG:NH2	33:BL:110:VAL:HG11	2.30	0.46
34:BM:8:LYS:CE	34:BM:70:ASP:HA	2.44	0.46
36:BO:9:ARG:HG3	36:BO:12:THR:HG21	1.97	0.46
36:BO:16:ARG:CD	36:BO:20:GLU:HG3	2.45	0.46
39:BR:5:PHE:CG	39:BR:6:GLN:N	2.84	0.46
40:BS:103:ILE:HG23	40:BS:104:THR:N	2.29	0.46
41:BT:12:ARG:HH21	44:BX:29:ARG:CZ	2.28	0.46
1:CA:1342:C:H4'	8:CI:126:PHE:O	2.15	0.46
1:CA:484:G:O2'	1:CA:485:U:OP2	2.28	0.46
1:CA:49:U:O2'	1:CA:50:A:H2'	2.14	0.46
1:CA:60:A:C4'	1:CA:61:G:O5'	2.63	0.46
1:CA:633:G:H2'	1:CA:634:C:H6	1.81	0.46
1:CA:699:C:C3'	1:CA:700:G:H5''	2.44	0.46
1:CA:754:C:H3'	1:CA:754:C:O2	2.15	0.46
1:CA:836:G:O2'	1:CA:837:U:H5'	2.15	0.46
20:CB:63:LYS:HD2	20:CB:87:ASP:OD2	2.14	0.46
2:CC:86:LEU:O	2:CC:89:VAL:HG22	2.14	0.46
4:CE:43:GLY:O	4:CE:72:ASN:HA	2.15	0.46
7:CH:74:ILE:O	7:CH:74:ILE:HG23	2.15	0.46
8:CI:108:ARG:HG3	8:CI:108:ARG:O	2.15	0.46
10:CK:110:THR:HB	21:CU:3:ILE:O	2.15	0.46
12:CM:47:LEU:HD12	12:CM:51:GLN:HB2	1.98	0.46
14:CO:48:ASP:CG	14:CO:51:SER:HB2	2.35	0.46
50:D3:12:ARG:CD	50:D3:23:HIS:HB2	2.45	0.46
23:DB:1328:A:H2'	23:DB:1330:C:C4	2.50	0.46
23:DB:131:A:H2'	23:DB:132:G:C8	2.48	0.46
23:DB:1450:G:N2	23:DB:1452:G:H1	1.95	0.46
23:DB:1526:C:H2'	23:DB:1527:G:O4'	2.14	0.46
23:DB:1707:G:H2'	23:DB:1708:C:H6	1.80	0.46
23:DB:1947:C:O2'	23:DB:1948:G:H5'	2.15	0.46
23:DB:2025:C:H2'	23:DB:2026:U:H6	1.77	0.46
23:DB:2299:U:H2'	23:DB:2300:C:H6	1.80	0.46
23:DB:2529:G:C4'	29:DG:175:LYS:HD3	2.46	0.46
23:DB:27:G:H1'	23:DB:513:A:N6	2.30	0.46
23:DB:2800:A:C4	23:DB:2801:G:H1'	2.51	0.46
23:DB:309:A:N3	23:DB:329:G:O2'	2.42	0.46
23:DB:39:G:H2'	23:DB:40:U:H6	1.80	0.46
25:DC:243:PRO:CB	25:DC:248:GLY:HA2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:58:LYS:HG3	25:DC:58:LYS:O	2.15	0.46
25:DC:28:PRO:CG	25:DC:79:ARG:NE	2.76	0.46
26:DD:118:PHE:HA	26:DD:164:GLN:HG2	1.96	0.46
27:DE:143:LEU:HD13	27:DE:185:LYS:HZ2	1.80	0.46
27:DE:161:ALA:HB1	27:DE:168:ASP:O	2.15	0.46
28:DF:69:ALA:HB3	28:DF:81:GLY:H	1.81	0.46
52:DI:16:MET:N	52:DI:42:ASN:OD1	2.49	0.46
27:DE:184:ASP:HB3	33:DL:13:LYS:HE3	1.97	0.46
23:DB:588:U:H5'	33:DL:29:LYS:HZ2	1.81	0.46
33:DL:63:LYS:HB2	50:D3:26:ALA:CB	2.43	0.46
34:DM:62:LYS:N	34:DM:104:GLU:HB3	2.20	0.46
36:DO:31:THR:HG22	36:DO:33:ARG:H	1.79	0.46
23:DB:533:G:N2	38:DQ:40:LYS:HD2	2.30	0.46
38:DQ:39:ILE:HG13	38:DQ:40:LYS:H	1.79	0.46
45:DY:43:ILE:HD13	45:DY:43:ILE:C	2.35	0.46
1:AA:1201:A:C1'	1:AA:1202:U:OP2	2.58	0.46
1:AA:1332:A:H2'	1:AA:1333:A:H8	1.79	0.46
1:AA:1382:C:H4'	6:AG:78:ARG:HD2	1.97	0.46
20:AB:100:LEU:O	20:AB:178:LEU:HG	2.16	0.46
20:AB:95:TRP:CH2	20:AB:100:LEU:HB2	2.50	0.46
2:AC:39:ARG:NH1	13:AN:91:GLU:HG2	2.31	0.46
2:AC:59:PRO:HD2	2:AC:62:SER:O	2.15	0.46
1:AA:619:U:O2	3:AD:129:VAL:HG13	2.15	0.46
3:AD:24:VAL:O	3:AD:25:ARG:C	2.53	0.46
3:AD:2:ARG:O	3:AD:3:TYR:HB2	2.15	0.46
5:AF:6:ILE:HD12	5:AF:88:MET:O	2.15	0.46
6:AG:125:ASP:HB3	6:AG:130:LYS:CB	2.35	0.46
8:AI:112:ARG:CB	8:AI:112:ARG:HH11	2.27	0.46
8:AI:40:ARG:HA	8:AI:44:ARG:HH21	1.81	0.46
1:AA:1180:A:P	8:AI:98:ARG:HH22	2.38	0.46
10:AK:41:LEU:HB3	10:AK:76:TYR:HE1	1.81	0.46
10:AK:56:LYS:O	10:AK:58:THR:N	2.49	0.46
11:AL:79:ILE:HD12	11:AL:79:ILE:C	2.35	0.46
11:AL:86:VAL:HG11	11:AL:89:LEU:HD23	1.97	0.46
13:AN:45:LEU:C	13:AN:45:LEU:HD23	2.36	0.46
14:AO:70:LYS:HZ1	14:AO:74:VAL:CG1	2.28	0.46
16:AQ:10:ARG:O	16:AQ:22:VAL:HG13	2.15	0.46
21:AU:20:ARG:CA	21:AU:24:LYS:HG3	2.39	0.46
49:B2:30:VAL:O	49:B2:33:ARG:HG3	2.15	0.46
33:BL:63:LYS:HB2	50:B3:11:LYS:HZ2	1.80	0.46
23:BB:1154:G:OP2	38:BQ:57:ARG:NH1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1623:G:O2'	23:BB:1624:U:H5'	2.15	0.46
23:BB:1666:G:O2'	23:BB:1667:G:H5'	2.14	0.46
23:BB:2699:C:O2'	23:BB:2700:A:H5'	2.16	0.46
23:BB:425:G:H2'	23:BB:426:C:C6	2.50	0.46
23:BB:857:G:H5'	43:BW:55:ASP:OD2	2.15	0.46
25:BC:155:ARG:O	25:BC:157:ALA:N	2.48	0.46
25:BC:243:PRO:HB2	25:BC:248:GLY:H	1.81	0.46
27:BE:177:PRO:HA	27:BE:180:LEU:HD12	1.95	0.46
31:BJ:45:THR:HB	38:BQ:63:ARG:NH1	2.30	0.46
34:BM:124:LEU:HD23	34:BM:124:LEU:N	2.31	0.46
24:BV:82:TYR:CD1	34:BM:36:VAL:HB	2.51	0.46
34:BM:82:MET:HB2	34:BM:84:LYS:CE	2.45	0.46
41:BT:28:ASN:ND2	41:BT:28:ASN:N	2.62	0.46
46:BZ:59:ARG:HH21	46:BZ:62:LYS:HZ2	1.61	0.46
1:CA:1434:A:H2'	1:CA:1435:G:C8	2.50	0.46
1:CA:336:A:O2'	1:CA:337:G:H5'	2.15	0.46
1:CA:546:A:OP1	3:CD:69:ARG:HB2	2.15	0.46
1:CA:691:G:H2'	1:CA:692:U:C6	2.51	0.46
1:CA:1190:G:P	2:CC:4:VAL:HG12	2.56	0.46
1:CA:409:U:OP1	3:CD:23:GLY:HA3	2.15	0.46
6:CG:78:ARG:HH11	6:CG:81:GLY:N	2.12	0.46
11:CL:30:ARG:HG2	11:CL:31:GLY:N	2.30	0.46
12:CM:106:ARG:CA	12:CM:106:ARG:HH11	2.28	0.46
19:CT:69:ASN:H	19:CT:69:ASN:HD22	1.60	0.46
48:D1:24:LYS:NZ	48:D1:24:LYS:HB2	2.31	0.46
51:D4:15:LYS:HB2	51:D4:15:LYS:HE3	1.79	0.46
23:DB:138:U:H2'	23:DB:140:C:N1	2.30	0.46
23:DB:1940:U:H1'	23:DB:1941:C:H5	1.81	0.46
23:DB:2616:C:H2'	23:DB:2617:U:H6	1.81	0.46
23:DB:2648:G:H2'	23:DB:2649:C:H6	1.80	0.46
23:DB:2699:C:O2'	23:DB:2700:A:H5'	2.15	0.46
23:DB:870:U:O2'	23:DB:871:U:H5'	2.15	0.46
23:DB:988:A:OP1	45:DY:10:ARG:HB3	2.15	0.46
52:DI:78:LEU:HA	52:DI:81:LYS:HE2	1.98	0.46
55:DB:3213:HOH:O	33:DL:41:ARG:HB2	2.14	0.46
34:DM:14:LYS:HB3	34:DM:72:PRO:HG3	1.97	0.46
38:DQ:52:ARG:CG	38:DQ:53:LYS:N	2.79	0.46
43:DW:82:GLU:HG3	43:DW:83:ALA:N	2.29	0.46
46:DZ:54:GLY:O	46:DZ:57:VAL:HB	2.15	0.46
1:AA:1016:A:H5'	1:AA:1218:C:H4'	1.98	0.46
1:AA:1149:C:O2'	1:AA:1150:A:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1181:G:OP2	1:AA:1181:G:O4'	2.33	0.46
1:AA:1462:C:H2'	1:AA:1463:U:C6	2.50	0.46
1:AA:512:U:H2'	1:AA:513:C:C6	2.49	0.46
1:AA:633:G:H2'	1:AA:634:C:H6	1.80	0.46
1:AA:812:G:H4'	1:AA:812:G:OP1	2.16	0.46
5:AF:32:ALA:HB1	5:AF:70:VAL:HG11	1.97	0.46
6:AG:41:ILE:CD1	6:AG:116:ALA:HB2	2.46	0.46
10:AK:80:ASN:OD1	10:AK:105:ARG:HB3	2.15	0.46
10:AK:125:LYS:O	10:AK:126:ARG:O	2.33	0.46
48:B1:10:LEU:HD22	48:B1:25:ASN:ND2	2.31	0.46
23:BB:1368:G:H5'	49:B2:25:LYS:HD3	1.98	0.46
33:BL:59:ARG:NH1	50:B3:51:LYS:HB2	2.31	0.46
22:BA:18:G:H2'	22:BA:19:C:H6	1.81	0.46
23:BB:1683:U:H2'	23:BB:1684:G:H8	1.80	0.46
23:BB:169:G:O2'	23:BB:170:U:H5'	2.15	0.46
23:BB:1824:G:H2'	23:BB:1825:U:C6	2.50	0.46
23:BB:1791:A:N1	23:BB:1829:A:H5'	2.30	0.46
23:BB:2515:C:H2'	23:BB:2516:A:H8	1.81	0.46
23:BB:2637:U:C2'	23:BB:2638:G:H5'	2.45	0.46
23:BB:2716:C:O2'	23:BB:2717:C:H5'	2.16	0.46
23:BB:692:C:H2'	23:BB:693:A:H8	1.80	0.46
23:BB:811:U:OP2	33:BL:31:GLY:HA2	2.16	0.46
25:BC:141:HIS:HB3	25:BC:190:THR:HB	1.97	0.46
25:BC:257:ARG:O	25:BC:261:ARG:NH1	2.49	0.46
27:BE:102:ARG:O	27:BE:105:LEU:HG	2.15	0.46
27:BE:27:LEU:HD21	27:BE:100:MET:HG2	1.96	0.46
28:BF:165:GLY:HA2	28:BF:168:LEU:CD2	2.45	0.46
30:BH:75:LEU:HD23	30:BH:76:GLU:N	2.31	0.46
30:BH:89:LYS:HE2	30:BH:89:LYS:CA	2.46	0.46
33:BL:120:VAL:HG13	33:BL:120:VAL:O	2.16	0.46
36:BO:19:GLN:HE21	43:BW:43:LYS:HE3	1.80	0.46
23:BB:1252:G:H22	38:BQ:36:GLN:NE2	2.14	0.46
39:BR:82:HIS:ND1	39:BR:82:HIS:O	2.49	0.46
42:BU:29:SER:O	42:BU:30:SER:HB2	2.16	0.46
43:BW:75:ASN:HB3	43:BW:76:ARG:H	1.62	0.46
46:BZ:35:ASP:C	46:BZ:42:PRO:HB3	2.35	0.46
46:BZ:5:ILE:CG2	46:BZ:63:ARG:HE	2.27	0.46
46:BZ:58:ASP:CG	46:BZ:63:ARG:HH11	2.19	0.46
1:CA:1349:A:H2'	1:CA:1350:A:O4'	2.15	0.46
1:CA:1473:G:O2'	1:CA:1474:U:H5'	2.16	0.46
1:CA:148:G:N3	1:CA:1446:A:H2	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:691:G:H1'	1:CA:696:A:N6	2.30	0.46
1:CA:768:A:H5'	1:CA:1524:C:H1'	1.97	0.46
1:CA:812:G:OP1	1:CA:812:G:H4'	2.16	0.46
1:CA:824:G:O2'	1:CA:825:A:H5'	2.14	0.46
20:CB:218:ALA:O	20:CB:221:ARG:HG2	2.16	0.46
20:CB:48:MET:HB3	20:CB:198:VAL:O	2.15	0.46
2:CC:110:LEU:N	2:CC:110:LEU:HD23	2.29	0.46
2:CC:76:ILE:O	2:CC:83:VAL:HG23	2.15	0.46
4:CE:89:THR:C	4:CE:91:SER:N	2.69	0.46
5:CF:38:ARG:HH21	5:CF:63:ASN:CG	2.18	0.46
5:CF:92:THR:CG2	5:CF:93:LYS:H	2.28	0.46
6:CG:71:THR:HA	6:CG:95:ARG:HE	1.81	0.46
11:CL:86:VAL:HG22	11:CL:95:HIS:HE2	1.81	0.46
12:CM:10:ASP:HA	12:CM:44:ILE:CD1	2.46	0.46
12:CM:9:PRO:O	12:CM:44:ILE:HG12	2.15	0.46
12:CM:63:VAL:O	12:CM:68:LEU:HD12	2.15	0.46
13:CN:63:CYS:HB3	13:CN:68:ARG:N	2.23	0.46
16:CQ:64:ARG:HB3	16:CQ:64:ARG:NH1	2.30	0.46
48:D1:5:ARG:N	48:D1:27:ARG:HH22	2.12	0.46
23:DB:1042:G:H2'	23:DB:1043:C:H6	1.80	0.46
23:DB:1192:G:O2'	23:DB:1193:G:H5'	2.16	0.46
23:DB:1424:G:H2'	23:DB:1425:G:O4'	2.15	0.46
23:DB:1487:U:H2'	23:DB:1488:C:C6	2.51	0.46
23:DB:1623:G:O2'	23:DB:1624:U:H5'	2.15	0.46
23:DB:163:C:H2'	23:DB:164:C:O4'	2.15	0.46
23:DB:1885:A:H3'	23:DB:1886:U:C6	2.51	0.46
23:DB:2051:A:H4'	26:DD:145:SER:CB	2.44	0.46
23:DB:2070:A:H2'	23:DB:2071:A:H8	1.80	0.46
23:DB:231:A:H3'	23:DB:232:G:H8	1.81	0.46
23:DB:2714:G:H2'	23:DB:2715:C:H6	1.80	0.46
23:DB:322:A:C2	23:DB:340:A:C6	3.03	0.46
23:DB:692:C:H2'	23:DB:693:A:C8	2.51	0.46
25:DC:243:PRO:HB3	25:DC:248:GLY:HA2	1.98	0.46
25:DC:257:ARG:HG2	25:DC:257:ARG:NH1	2.31	0.46
26:DD:36:GLN:HG2	26:DD:88:GLU:HA	1.97	0.46
26:DD:62:LYS:HB2	26:DD:63:PRO:HD3	1.97	0.46
27:DE:116:ASP:CG	27:DE:185:LYS:HE2	2.36	0.46
28:DF:132:ARG:HD3	28:DF:133:GLU:N	2.25	0.46
29:DG:172:GLU:O	29:DG:173:ALA:C	2.54	0.46
29:DG:42:VAL:HB	29:DG:51:PHE:HA	1.96	0.46
31:DJ:106:LYS:C	31:DJ:108:MET:H	2.18	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:DL:79:LEU:HG	33:DL:111:ILE:O	2.15	0.46
34:DM:61:GLY:CA	34:DM:105:MET:HB2	2.45	0.46
34:DM:18:ARG:N	34:DM:18:ARG:CD	2.78	0.46
34:DM:73:ILE:HD11	34:DM:92:TRP:HE3	1.80	0.46
34:DM:75:GLU:O	34:DM:86:LYS:HB3	2.15	0.46
35:DN:4:ARG:HB3	35:DN:4:ARG:NH1	2.30	0.46
22:DA:38:C:H4'	36:DO:100:HIS:CE1	2.51	0.46
37:DP:13:LYS:HZ1	37:DP:15:ASP:HB2	1.79	0.46
39:DR:4:VAL:O	39:DR:41:ILE:HG12	2.15	0.46
23:DB:2091:C:O4'	46:DZ:49:ARG:NE	2.49	0.46
1:AA:1220:G:H4'	18:AS:33:TRP:O	2.16	0.46
1:AA:529:G:H22	11:AL:47:ALA:HB2	1.80	0.46
1:AA:834:U:H2'	1:AA:835:U:C6	2.50	0.46
1:AA:840:C:N3	1:AA:842:U:H4'	2.30	0.46
20:AB:187:ASP:CG	20:AB:188:THR:N	2.69	0.46
3:AD:54:LEU:HA	3:AD:202:LEU:HD22	1.97	0.46
4:AE:45:VAL:HG12	4:AE:116:VAL:HG23	1.98	0.46
4:AE:41:GLY:HA2	4:AE:118:GLY:HA2	1.97	0.46
7:AH:51:GLU:O	7:AH:56:PRO:HA	2.15	0.46
8:AI:46:VAL:HA	8:AI:49:GLN:HB2	1.98	0.46
9:AJ:30:LYS:HG3	9:AJ:36:VAL:HB	1.98	0.46
11:AL:27:PRO:HG2	11:AL:28:GLN:NE2	2.31	0.46
50:B3:4:LYS:HE3	50:B3:58:ILE:HB	1.98	0.46
51:B4:24:ARG:HD3	51:B4:26:ILE:HG23	1.98	0.46
22:BA:117:G:OP1	36:BO:55:GLU:OE1	2.33	0.46
23:BB:1453:A:N7	35:BN:72:ASP:OD1	2.49	0.46
23:BB:165:A:H2'	23:BB:166:U:H6	1.80	0.46
23:BB:1795:C:H2'	23:BB:1796:U:C6	2.51	0.46
23:BB:2379:G:H2'	23:BB:2380:C:C6	2.51	0.46
23:BB:2835:A:H61	23:BB:2878:U:H2'	1.81	0.46
23:BB:322:A:C2	23:BB:340:A:C6	3.03	0.46
23:BB:459:U:H5'	49:B2:41:ARG:HH21	1.80	0.46
23:BB:1820:U:H3	25:BC:197:ALA:HA	1.79	0.46
25:BC:45:ASN:O	25:BC:46:GLY:C	2.54	0.46
25:BC:59:GLN:O	25:BC:60:ALA:HB3	2.15	0.46
25:BC:37:SER:H	25:BC:62:ARG:CB	2.29	0.46
25:BC:79:ARG:HH11	25:BC:80:LEU:N	2.11	0.46
26:BD:114:LYS:HE3	26:BD:196:ALA:CB	2.45	0.46
30:BH:3:VAL:CG2	30:BH:21:VAL:HG21	2.45	0.46
52:BI:125:THR:O	52:BI:129:GLU:HG3	2.16	0.46
52:BI:102:ARG:HB2	52:BI:141:ASP:OD2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:BL:63:LYS:CA	50:B3:24:LYS:HD3	2.46	0.46
33:BL:83:ALA:O	33:BL:84:LYS:O	2.34	0.46
34:BM:66:ARG:H	34:BM:100:LYS:HB2	1.80	0.46
34:BM:108:VAL:HG23	34:BM:112:LEU:CD1	2.40	0.46
35:BN:73:ASN:CA	35:BN:76:VAL:HG13	2.45	0.46
36:BO:26:LEU:HD13	36:BO:92:PHE:O	2.16	0.46
38:BQ:97:ILE:HA	38:BQ:100:PHE:CE2	2.51	0.46
40:BS:34:ASP:CA	40:BS:37:THR:HG22	2.45	0.46
40:BS:69:LEU:O	40:BS:110:ARG:HD2	2.15	0.46
41:BT:6:ARG:C	41:BT:8:LEU:H	2.18	0.46
24:BV:44:HIS:O	24:BV:46:LYS:N	2.49	0.46
46:BZ:12:ILE:HG22	46:BZ:24:ILE:HD11	1.96	0.46
1:CA:109:A:H4'	1:CA:110:C:OP2	2.16	0.46
1:CA:1131:G:H2'	1:CA:1132:C:H5'	1.98	0.46
1:CA:1411:C:O2'	1:CA:1412:C:H5'	2.15	0.46
1:CA:1432:G:H1'	1:CA:1468:A:N6	2.31	0.46
1:CA:1492:A:N6	1:CA:1494:G:N9	2.64	0.46
1:CA:1505:G:N7	53:CA:1601:KSG:H3	2.31	0.46
1:CA:455:G:H2'	1:CA:456:A:C8	2.51	0.46
1:CA:782:A:H4'	1:CA:1514:G:O2'	2.16	0.46
20:CB:113:LEU:HD23	20:CB:113:LEU:C	2.36	0.46
2:CC:174:LEU:HD21	2:CC:200:TRP:CD1	2.51	0.46
2:CC:55:VAL:HG23	2:CC:66:THR:HB	1.98	0.46
4:CE:95:MET:HG3	4:CE:124:ALA:CB	2.44	0.46
6:CG:8:GLN:HA	6:CG:8:GLN:HE21	1.81	0.46
8:CI:8:THR:OG1	8:CI:9:GLY:N	2.49	0.46
13:CN:30:ILE:HB	13:CN:44:VAL:CB	2.44	0.46
13:CN:76:PHE:O	13:CN:78:LEU:HD13	2.15	0.46
16:CQ:63:CYS:SG	16:CQ:66:LEU:HD11	2.56	0.46
48:D1:27:ARG:O	48:D1:27:ARG:HG2	2.15	0.46
49:D2:19:ARG:HD3	49:D2:19:ARG:C	2.35	0.46
23:DB:1750:G:H2'	23:DB:1751:U:C6	2.50	0.46
23:DB:1827:U:O2'	23:DB:1828:G:H5'	2.14	0.46
23:DB:1841:U:H2'	23:DB:1842:G:H8	1.79	0.46
23:DB:2581:G:OP1	26:DD:134:HIS:CD2	2.69	0.46
23:DB:458:G:C2'	23:DB:459:U:OP2	2.64	0.46
23:DB:934:U:H2'	23:DB:935:C:H6	1.80	0.46
23:DB:948:C:H2'	23:DB:949:G:H8	1.81	0.46
25:DC:51:ARG:NH1	25:DC:54:GLY:HA3	2.20	0.46
26:DD:21:SER:O	26:DD:23:PRO:HD3	2.16	0.46
27:DE:116:ASP:C	27:DE:117:ARG:HD2	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:148:ILE:HA	27:DE:185:LYS:CB	2.35	0.46
27:DE:116:ASP:OD2	27:DE:185:LYS:HG3	2.16	0.46
30:DH:124:THR:HG23	30:DH:128:HIS:CE1	2.49	0.46
30:DH:3:VAL:HG22	30:DH:21:VAL:CG1	2.28	0.46
38:DQ:39:ILE:CG1	38:DQ:40:LYS:N	2.79	0.46
39:DR:41:ILE:HD12	39:DR:42:ALA:H	1.80	0.46
39:DR:64:VAL:HG22	39:DR:65:ALA:N	2.23	0.46
24:DV:93:ARG:HG2	24:DV:94:ALA:N	2.30	0.46
44:DX:31:GLN:O	44:DX:32:ALA:HB3	2.15	0.46
44:DX:28:LEU:HD13	44:DX:42:LEU:CG	2.45	0.46
1:AA:1010:U:H2'	1:AA:1011:C:C6	2.50	0.46
1:AA:1310:G:O2'	1:AA:1311:A:H5'	2.15	0.46
1:AA:1480:A:O2'	1:AA:1481:U:H5'	2.16	0.46
1:AA:328:C:H4'	1:AA:329:A:H5''	1.98	0.46
1:AA:334:C:H2'	1:AA:335:C:H6	1.80	0.46
1:AA:807:A:H2'	1:AA:808:C:C6	2.51	0.46
1:AA:836:G:O2'	1:AA:837:U:H5'	2.15	0.46
1:AA:920:U:H2'	1:AA:921:U:H6	1.81	0.46
3:AD:131:ILE:O	3:AD:134:TYR:HB2	2.16	0.46
3:AD:24:VAL:O	3:AD:25:ARG:HD2	2.15	0.46
4:AE:17:VAL:HA	4:AE:34:ALA:HA	1.98	0.46
7:AH:93:LYS:HD3	7:AH:93:LYS:N	2.30	0.46
9:AJ:39:PRO:HA	9:AJ:74:VAL:HG22	1.98	0.46
11:AL:71:HIS:ND1	11:AL:72:ASN:N	2.63	0.46
12:AM:1:ALA:O	12:AM:8:ILE:HG22	2.15	0.46
23:BB:1021:A:N6	23:BB:1141:U:H3	1.98	0.46
23:BB:1353:A:H2'	23:BB:1354:A:C8	2.50	0.46
23:BB:144:A:O2'	23:BB:145:C:H5'	2.16	0.46
23:BB:2031:A:C6	23:BB:2498:C:H1'	2.50	0.46
23:BB:2063:C:O2	23:BB:2450:A:N1	2.49	0.46
23:BB:2393:U:OP1	50:B3:29:ARG:NE	2.49	0.46
23:BB:2648:G:H2'	23:BB:2649:C:H6	1.80	0.46
23:BB:2700:A:H2'	23:BB:2701:U:C6	2.51	0.46
23:BB:26:G:H2'	23:BB:27:G:O4'	2.15	0.46
23:BB:473:G:O2'	23:BB:474:G:H5'	2.15	0.46
23:BB:494:G:O2'	23:BB:495:G:H5'	2.15	0.46
23:BB:811:U:C4	33:BL:29:LYS:HD3	2.50	0.46
25:BC:127:ASN:ND2	25:BC:128:THR:N	2.55	0.46
25:BC:136:VAL:CG1	25:BC:165:ALA:HA	2.39	0.46
25:BC:224:MET:SD	25:BC:224:MET:N	2.88	0.46
25:BC:78:GLU:CB	25:BC:92:LEU:HB3	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:100:LEU:C	26:BD:102:ALA:H	2.19	0.46
26:BD:118:PHE:N	26:BD:164:GLN:HG3	2.31	0.46
26:BD:170:VAL:HG11	26:BD:194:PRO:CG	2.46	0.46
27:BE:119:ILE:HA	27:BE:189:THR:H	1.81	0.46
27:BE:153:LEU:H	27:BE:171:ASP:HB3	1.80	0.46
29:BG:32:LEU:HD21	29:BG:74:MET:CG	2.46	0.46
30:BH:78:VAL:CG1	30:BH:102:ALA:HB1	2.46	0.46
30:BH:78:VAL:HB	30:BH:143:ILE:O	2.16	0.46
31:BJ:81:ILE:CG1	31:BJ:82:GLY:N	2.69	0.46
33:BL:116:VAL:HG12	33:BL:117:THR:N	2.21	0.46
35:BN:28:LEU:HD21	35:BN:113:ILE:HG21	1.96	0.46
31:BJ:42:ALA:HB1	38:BQ:67:ALA:HB2	1.97	0.46
39:BR:39:LEU:HA	39:BR:60:LYS:O	2.15	0.46
42:BU:11:ILE:HB	42:BU:69:VAL:HG11	1.96	0.46
43:BW:56:HIS:HE1	43:BW:58:LEU:HB2	1.79	0.46
43:BW:24:ARG:HG3	43:BW:59:PHE:HD1	1.81	0.46
1:CA:1006:G:H2'	1:CA:1007:U:H6	1.80	0.46
1:CA:1349:A:OP1	8:CI:121:ARG:HB2	2.16	0.46
1:CA:1422:G:H5''	32:DK:48:PRO:CB	2.39	0.46
1:CA:599:C:O2'	1:CA:600:A:H5'	2.15	0.46
1:CA:628:G:H2'	1:CA:629:A:C8	2.51	0.46
1:CA:737:C:H5'	5:CF:89:VAL:O	2.15	0.46
4:CE:154:ALA:HB1	7:CH:65:PHE:CZ	2.50	0.46
8:CI:104:THR:HG22	8:CI:105:ARG:N	2.31	0.46
1:CA:1347:G:O5'	8:CI:108:ARG:HB2	2.16	0.46
8:CI:9:GLY:O	8:CI:16:ALA:HB3	2.15	0.46
10:CK:12:ARG:HD3	10:CK:13:LYS:NZ	2.29	0.46
10:CK:70:ALA:HB1	10:CK:74:LYS:HB2	1.97	0.46
12:CM:3:ILE:HA	12:CM:56:ARG:HH11	1.80	0.46
12:CM:79:LEU:HA	12:CM:82:LEU:CG	2.46	0.46
48:D1:47:ILE:O	48:D1:48:TYR:HB2	2.16	0.46
23:DB:1552:A:H2'	23:DB:1553:A:C5'	2.43	0.46
23:DB:1759:A:H4'	23:DB:2715:C:O4'	2.16	0.46
23:DB:1824:G:H2'	23:DB:1825:U:H6	1.80	0.46
23:DB:1919:A:H2'	23:DB:1920:C:H5'	1.97	0.46
23:DB:2208:C:H2'	23:DB:2209:G:C8	2.51	0.46
23:DB:2311:A:O2'	28:DF:39:VAL:HG21	2.15	0.46
23:DB:2386:A:H2'	23:DB:2387:U:H6	1.81	0.46
23:DB:2543:G:H2'	23:DB:2544:G:C8	2.50	0.46
23:DB:2645:G:H4'	23:DB:2732:G:H2'	1.96	0.46
23:DB:2815:C:C2	23:DB:2816:G:C8	3.04	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2816:G:N3	23:DB:2883:A:O2'	2.48	0.46
23:DB:307:G:N1	23:DB:310:A:OP2	2.48	0.46
23:DB:834:G:H2'	23:DB:835:C:H6	1.80	0.46
26:DD:146:ILE:N	26:DD:146:ILE:CD1	2.79	0.46
27:DE:23:PHE:C	27:DE:110:SER:HB2	2.36	0.46
27:DE:31:VAL:HG21	27:DE:104:ALA:CB	2.45	0.46
23:DB:1098:A:O3'	52:DI:4:VAL:O	2.33	0.46
52:DI:72:THR:OG1	52:DI:73:PRO:HD2	2.15	0.46
31:DJ:130:HIS:O	31:DJ:131:ASN:C	2.54	0.46
34:DM:96:ILE:HG22	34:DM:97:GLN:OE1	2.16	0.46
37:DP:91:VAL:HG23	37:DP:92:ARG:N	2.29	0.46
39:DR:63:VAL:HG13	39:DR:64:VAL:N	2.30	0.46
39:DR:81:LYS:O	39:DR:83:TYR:N	2.48	0.46
40:DS:23:LEU:HD13	40:DS:25:ARG:NH2	2.22	0.46
40:DS:23:LEU:HD23	47:D0:21:LEU:HD13	1.96	0.46
41:DT:32:LEU:O	41:DT:83:ALA:HB3	2.15	0.46
42:DU:42:LYS:H	42:DU:57:ILE:HG21	1.80	0.46
43:DW:54:ARG:HD2	43:DW:54:ARG:N	2.31	0.46
1:AA:1236:A:H2'	1:AA:1237:C:O4'	2.15	0.46
1:AA:239:U:H5'	1:AA:239:U:H6	1.80	0.46
1:AA:285:C:H2'	1:AA:286:C:H6	1.81	0.46
1:AA:81:A:C2	1:AA:82:G:N7	2.83	0.46
5:AF:29:ILE:HD13	5:AF:64:VAL:CG1	2.45	0.46
6:AG:94:ARG:CD	6:AG:98:LEU:HD11	2.43	0.46
8:AI:51:LEU:HB3	8:AI:56:MET:CG	2.45	0.46
10:AK:32:THR:HA	10:AK:43:TRP:HA	1.98	0.46
10:AK:88:PRO:HA	10:AK:92:ARG:HD2	1.98	0.46
11:AL:98:ARG:CB	11:AL:116:TYR:HA	2.45	0.46
18:AS:12:LEU:O	18:AS:15:LEU:HB3	2.15	0.46
23:BB:682:G:H5'	49:B2:26:ASN:OD1	2.16	0.46
51:B4:27:CYS:SG	51:B4:33:HIS:NE2	2.89	0.46
23:BB:1113:U:H5''	29:BG:2:ARG:H	1.79	0.46
23:BB:131:A:O2'	23:BB:132:G:H5'	2.15	0.46
23:BB:1747:U:H2'	23:BB:1748:C:H6	1.81	0.46
23:BB:215:G:H4'	23:BB:216:A:OP1	2.15	0.46
23:BB:2239:G:H5'	25:BC:246:PRO:HG2	1.97	0.46
23:BB:2297:A:N6	23:BB:2319:G:C5'	2.78	0.46
23:BB:248:G:C2	23:BB:2431:U:H4'	2.51	0.46
23:BB:247:G:H4'	23:BB:386:G:C5	2.51	0.46
23:BB:30:G:O2'	23:BB:31:C:H5'	2.16	0.46
23:BB:527:C:O2'	23:BB:528:A:C8	2.66	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:692:C:H2'	23:BB:693:A:C8	2.51	0.46
23:BB:1791:A:C4'	25:BC:207:ALA:H	2.28	0.46
25:BC:38:LYS:O	25:BC:39:SER:HB3	2.15	0.46
26:BD:22:ILE:HG12	26:BD:22:ILE:O	2.15	0.46
26:BD:38:LYS:N	26:BD:38:LYS:HD3	2.30	0.46
27:BE:35:TYR:CE2	27:BE:178:VAL:HG21	2.50	0.46
27:BE:28:VAL:HG23	27:BE:29:HIS:N	2.31	0.46
23:BB:615:U:O2	27:BE:38:GLY:HA3	2.15	0.46
29:BG:6:ALA:N	29:BG:7:PRO:CD	2.79	0.46
32:BK:23:LYS:CG	32:BK:24:VAL:H	2.27	0.46
32:BK:61:VAL:HG23	32:BK:85:VAL:HB	1.98	0.46
33:BL:28:GLY:O	33:BL:29:LYS:HB2	2.15	0.46
33:BL:88:GLY:O	33:BL:90:VAL:HG23	2.15	0.46
23:BB:2484:G:HO2'	34:BM:45:GLN:NE2	2.14	0.46
35:BN:9:GLN:HE21	35:BN:17:ARG:NH2	2.14	0.46
36:BO:15:ARG:HH22	36:BO:17:LYS:HD2	1.80	0.46
38:BQ:89:ILE:O	38:BQ:89:ILE:HG12	2.15	0.46
24:BV:4:ILE:HD11	24:BV:61:LEU:CD1	2.46	0.46
24:BV:77:VAL:HG21	24:BV:79:ARG:HH21	1.81	0.46
1:CA:1095:U:H2'	1:CA:1096:C:H6	1.81	0.46
1:CA:1388:C:H2'	1:CA:1389:C:H6	1.79	0.46
1:CA:1486:G:H2'	1:CA:1487:G:C8	2.51	0.46
1:CA:807:A:H2'	1:CA:808:C:C6	2.51	0.46
20:CB:19:THR:HA	20:CB:37:VAL:HB	1.98	0.46
2:CC:190:THR:HG22	2:CC:191:THR:H	1.80	0.46
2:CC:8:GLY:O	2:CC:11:LEU:HG	2.15	0.46
4:CE:87:VAL:O	4:CE:88:HIS:HB3	2.16	0.46
7:CH:38:VAL:CG2	7:CH:111:THR:HG22	2.46	0.46
10:CK:92:ARG:HH11	10:CK:92:ARG:HG2	1.81	0.46
11:CL:65:TYR:HB3	11:CL:95:HIS:CD2	2.51	0.46
13:CN:17:ASP:C	13:CN:19:TYR:H	2.19	0.46
13:CN:45:LEU:O	13:CN:48:GLN:HG3	2.16	0.46
15:CP:4:ILE:CD1	15:CP:57:ILE:HG12	2.44	0.46
19:CT:69:ASN:ND2	19:CT:69:ASN:N	2.53	0.46
21:CU:8:ASN:HB2	21:CU:9:GLU:OE2	2.16	0.46
49:D2:39:ARG:HG3	49:D2:39:ARG:HH11	1.79	0.46
23:DB:1019:U:O2'	23:DB:1020:A:H5'	2.15	0.46
23:DB:1108:U:H2'	23:DB:1109:C:H6	1.81	0.46
23:DB:1163:G:O2'	23:DB:1164:C:H5'	2.16	0.46
23:DB:1165:A:C2	23:DB:1166:G:N7	2.83	0.46
23:DB:1251:C:HO2'	23:DB:1252:G:H3'	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1537:G:H2'	23:DB:1538:G:O4'	2.14	0.46
23:DB:1795:C:H2'	23:DB:1796:U:C6	2.51	0.46
23:DB:2093:G:H2'	23:DB:2094:A:H8	1.81	0.46
23:DB:2156:G:H2'	23:DB:2157:G:C4'	2.37	0.46
23:DB:964:C:O2'	23:DB:2273:A:H1'	2.15	0.46
23:DB:2373:G:H2'	23:DB:2374:C:H6	1.80	0.46
23:DB:242:G:O2'	23:DB:243:U:OP2	2.33	0.46
23:DB:256:A:O2'	23:DB:257:C:H5'	2.16	0.46
23:DB:2603:G:O2'	23:DB:2604:U:H5'	2.16	0.46
23:DB:2722:G:H2'	23:DB:2723:C:H6	1.81	0.46
23:DB:277:G:H2'	23:DB:277:G:N3	2.31	0.46
23:DB:479:A:O2'	23:DB:480:A:OP2	2.30	0.46
23:DB:570:G:O2'	23:DB:571:U:H5'	2.16	0.46
25:DC:136:VAL:C	25:DC:138:SER:N	2.69	0.46
27:DE:188:MET:CE	27:DE:190:ALA:HB2	2.46	0.46
29:DG:42:VAL:HA	29:DG:50:THR:O	2.16	0.46
29:DG:94:ARG:H	29:DG:105:SER:HB2	1.80	0.46
30:DH:126:GLY:O	30:DH:145:ASN:HA	2.16	0.46
52:DI:18:ASN:N	52:DI:19:PRO:CD	2.78	0.46
32:DK:11:ALA:C	32:DK:100:PHE:H	2.20	0.46
32:DK:4:GLU:OE2	32:DK:23:LYS:HD2	2.15	0.46
32:DK:84:CYS:O	32:DK:85:VAL:HB	2.16	0.46
33:DL:92:LEU:HB2	33:DL:125:LEU:CD1	2.45	0.46
23:DB:2415:G:H4'	33:DL:66:PHE:CB	2.46	0.46
35:DN:76:VAL:O	35:DN:80:PHE:HD1	1.98	0.46
36:DO:100:HIS:HA	36:DO:104:GLN:NE2	2.30	0.46
37:DP:52:ARG:HG2	37:DP:60:VAL:CG2	2.46	0.46
38:DQ:52:ARG:HB3	38:DQ:52:ARG:HH11	1.79	0.46
38:DQ:87:VAL:O	39:DR:54:VAL:HG21	2.16	0.46
39:DR:79:ARG:HB3	39:DR:89:HIS:HE1	1.81	0.46
42:DU:42:LYS:N	42:DU:57:ILE:HG21	2.31	0.46
42:DU:38:ILE:HG13	42:DU:62:ALA:CB	2.45	0.46
43:DW:44:PHE:CD2	43:DW:77:LYS:HB3	2.50	0.46
43:DW:57:THR:HG22	43:DW:77:LYS:HG2	1.98	0.46
45:DY:26:LEU:C	45:DY:28:LEU:H	2.19	0.46
1:AA:1202:U:H2'	1:AA:1203:C:C5'	2.46	0.46
1:AA:28:A:O2'	1:AA:296:U:OP1	2.32	0.46
1:AA:428:G:H4'	1:AA:429:U:O5'	2.16	0.46
1:AA:628:G:H2'	1:AA:629:A:C8	2.51	0.46
1:AA:820:U:H4'	1:AA:821:G:OP2	2.16	0.46
20:AB:15:PHE:O	20:AB:40:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AB:164:ASP:CG	20:AB:203:ASP:HB2	2.36	0.46
2:AC:111:ASP:O	2:AC:115:VAL:HG23	2.15	0.46
2:AC:38:VAL:HG23	2:AC:39:ARG:N	2.29	0.46
6:AG:119:LEU:HD13	6:AG:119:LEU:C	2.36	0.46
1:AA:1372:U:OP1	8:AI:72:SER:HB2	2.16	0.46
9:AJ:15:HIS:CD2	9:AJ:19:ASP:HB2	2.51	0.46
12:AM:58:GLU:HA	12:AM:61:LYS:HE2	1.98	0.46
19:AT:60:GLN:CD	19:AT:60:GLN:H	2.19	0.46
21:AU:15:LEU:O	21:AU:18:PHE:HB2	2.16	0.46
47:B0:2:VAL:CG2	47:B0:3:GLN:N	2.79	0.46
48:B1:10:LEU:HD22	48:B1:25:ASN:HD22	1.81	0.46
23:BB:1083:U:C2	23:BB:1086:A:N1	2.83	0.46
23:BB:1091:G:O2'	23:BB:1092:C:H5'	2.15	0.46
23:BB:1712:U:H3'	23:BB:1713:A:C5'	2.46	0.46
23:BB:1735:A:H2'	23:BB:1736:U:C6	2.50	0.46
23:BB:1936:A:H2	23:BB:1943:U:C5	2.33	0.46
23:BB:1946:U:C2	23:BB:1947:C:C5	3.03	0.46
23:BB:2460:U:H2'	23:BB:2461:A:H8	1.81	0.46
23:BB:2714:G:H2'	23:BB:2715:C:C6	2.51	0.46
23:BB:2723:C:C5'	35:BN:3:HIS:HB2	2.46	0.46
23:BB:2815:C:C2	23:BB:2816:G:C8	3.04	0.46
23:BB:2816:G:N3	23:BB:2883:A:O2'	2.48	0.46
23:BB:774:G:OP2	25:BC:47:ARG:HD3	2.15	0.46
23:BB:788:A:H1'	49:B2:4:THR:OG1	2.15	0.46
25:BC:115:ILE:HG23	25:BC:126:GLY:O	2.15	0.46
25:BC:62:ARG:O	25:BC:62:ARG:NE	2.48	0.46
25:BC:27:LYS:HG3	25:BC:80:LEU:HB2	1.96	0.46
26:BD:131:ASP:CG	26:BD:132:ALA:N	2.69	0.46
26:BD:135:GLY:C	26:BD:137:SER:N	2.69	0.46
26:BD:148:GLN:HB3	26:BD:151:THR:CG2	2.45	0.46
26:BD:23:PRO:HA	26:BD:190:LYS:HA	1.98	0.46
26:BD:1:MET:CB	26:BD:84:LEU:HD22	2.40	0.46
28:BF:35:LEU:O	28:BF:36:ASN:HB3	2.14	0.46
29:BG:94:ARG:NH1	29:BG:127:GLN:HB2	2.31	0.46
30:BH:3:VAL:HB	30:BH:37:VAL:CG2	2.46	0.46
52:BI:46:ASP:HA	52:BI:50:LYS:HG3	1.96	0.46
31:BJ:17:VAL:HG13	31:BJ:57:LEU:HD21	1.97	0.46
31:BJ:19:ASP:O	31:BJ:20:ALA:HB3	2.16	0.46
32:BK:99:ILE:CG2	32:BK:119:ALA:HB2	2.46	0.46
33:BL:120:VAL:H	33:BL:138:ALA:CB	2.29	0.46
35:BN:41:ALA:HB1	35:BN:97:ILE:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:78:LYS:O	35:BN:82:GLU:HB2	2.15	0.46
36:BO:64:TYR:CD1	36:BO:74:VAL:HG21	2.50	0.46
36:BO:84:GLU:N	36:BO:84:GLU:OE1	2.49	0.46
37:BP:57:ALA:C	37:BP:59:THR:H	2.19	0.46
37:BP:9:GLN:CD	37:BP:56:SER:HB2	2.36	0.46
26:BD:11:MET:HE1	37:BP:9:GLN:HG2	1.98	0.46
38:BQ:5:ARG:CD	38:BQ:5:ARG:H	2.27	0.46
23:BB:1011:G:H5''	38:BQ:76:SER:OG	2.16	0.46
44:BX:30:MET:CG	44:BX:31:GLN:N	2.79	0.46
44:BX:59:GLU:C	44:BX:61:ALA:H	2.19	0.46
23:BB:396:G:C4'	46:BZ:28:VAL:HG21	2.34	0.46
1:CA:278:G:H21	1:CA:279:A:N6	2.10	0.46
1:CA:601:G:H2'	1:CA:602:A:H8	1.80	0.46
1:CA:684:U:O2'	10:CK:39:ASN:HB3	2.16	0.46
1:CA:721:G:H5''	17:CR:42:ARG:HH21	1.81	0.46
2:CC:177:LEU:HD22	2:CC:177:LEU:N	2.31	0.46
4:CE:76:ASN:O	4:CE:77:ASN:HB3	2.16	0.46
8:CI:117:LEU:CD2	8:CI:123:ARG:HD3	2.46	0.46
8:CI:87:MET:HB2	8:CI:94:ARG:NH2	2.31	0.46
11:CL:23:LEU:O	11:CL:25:ALA:N	2.49	0.46
18:CS:43:MET:O	18:CS:46:LEU:HB2	2.16	0.46
49:D2:43:THR:C	49:D2:44:VAL:HG22	2.36	0.46
23:DB:1309:G:C5'	49:D2:7:PRO:HB2	2.46	0.46
22:DA:75:G:N1	22:DA:102:G:N2	2.64	0.46
23:DB:1162:G:O2'	23:DB:1163:G:H5'	2.16	0.46
23:DB:139:U:H3'	23:DB:139:U:OP2	2.16	0.46
23:DB:153:U:H2'	23:DB:154:U:O4'	2.16	0.46
23:DB:2083:G:H2'	23:DB:2084:C:C6	2.51	0.46
23:DB:267:C:H2'	23:DB:268:C:C6	2.50	0.46
23:DB:2716:C:O2'	23:DB:2717:C:H5'	2.15	0.46
23:DB:2721:A:H2'	23:DB:2722:G:C8	2.51	0.46
23:DB:2800:A:N3	23:DB:2801:G:H1'	2.30	0.46
23:DB:1791:A:OP1	25:DC:211:ARG:HG2	2.16	0.46
25:DC:38:LYS:HG3	25:DC:39:SER:N	2.31	0.46
25:DC:92:LEU:HD12	25:DC:101:ARG:O	2.16	0.46
26:DD:185:ASN:O	26:DD:186:LEU:HD12	2.16	0.46
26:DD:29:VAL:CG2	26:DD:30:GLU:H	2.18	0.46
27:DE:154:ASP:OD2	27:DE:156:ASN:HB3	2.16	0.46
23:DB:443:A:N7	27:DE:40:ARG:HG2	2.31	0.46
52:DI:54:ILE:O	52:DI:54:ILE:HG23	2.16	0.46
31:DJ:17:VAL:HA	31:DJ:55:ILE:HB	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DK:43:ILE:HG22	32:DK:54:LYS:HA	1.97	0.46
32:DK:18:ARG:HB2	32:DK:45:GLU:HG2	1.98	0.46
33:DL:118:THR:O	33:DL:120:VAL:N	2.49	0.46
33:DL:62:PRO:HG3	50:D3:12:ARG:NH2	2.29	0.46
36:DO:82:ALA:HB1	36:DO:87:ILE:CG2	2.46	0.46
38:DQ:108:LEU:HA	38:DQ:111:LYS:CD	2.45	0.46
38:DQ:39:ILE:HA	38:DQ:43:GLN:NE2	2.30	0.46
38:DQ:53:LYS:N	38:DQ:53:LYS:HE2	2.13	0.46
39:DR:35:PHE:HB3	39:DR:64:VAL:HG12	1.97	0.46
23:DB:814:C:OP1	39:DR:87:GLN:O	2.34	0.46
41:DT:35:ALA:O	41:DT:81:LYS:HD3	2.16	0.46
43:DW:38:ARG:O	43:DW:39:GLN:HB3	2.15	0.46
1:AA:1096:C:H2'	1:AA:1097:C:H6	1.81	0.46
1:AA:1304:G:H2'	1:AA:1305:G:C1'	2.46	0.46
1:AA:132:C:O2'	1:AA:133:U:H5'	2.15	0.46
1:AA:1427:C:O2'	1:AA:1428:A:H5'	2.16	0.46
1:AA:1460:C:O2'	1:AA:1461:G:H5'	2.16	0.46
1:AA:382:A:O2'	1:AA:383:A:H5'	2.15	0.46
1:AA:41:G:H2'	1:AA:42:G:H8	1.81	0.46
1:AA:599:C:O2'	1:AA:600:A:H5'	2.15	0.46
1:AA:601:G:H2'	1:AA:602:A:C8	2.50	0.46
1:AA:82:G:C6	1:AA:88:U:O2	2.69	0.46
2:AC:173:PRO:C	2:AC:175:HIS:H	2.19	0.46
3:AD:98:ASP:CB	3:AD:132:ALA:HB1	2.46	0.46
4:AE:82:HIS:HB2	4:AE:83:PRO:HD2	1.98	0.46
5:AF:50:PRO:HA	5:AF:54:LEU:O	2.16	0.46
6:AG:11:ILE:HG12	6:AG:24:LYS:HE2	1.97	0.46
7:AH:86:LYS:HB3	7:AH:90:GLU:HB2	1.97	0.46
8:AI:119:LYS:NZ	8:AI:119:LYS:HB3	2.31	0.46
10:AK:63:GLN:O	10:AK:67:GLU:HB3	2.15	0.46
11:AL:120:ARG:HG3	11:AL:120:ARG:HH11	1.81	0.46
11:AL:35:ARG:O	11:AL:52:CYS:HB2	2.16	0.46
15:AP:6:LEU:HD11	15:AP:71:VAL:HB	1.98	0.46
18:AS:12:LEU:HD23	18:AS:12:LEU:O	2.15	0.46
18:AS:29:PRO:HB3	18:AS:47:THR:CG2	2.46	0.46
19:AT:60:GLN:HE21	19:AT:61:ALA:H	1.64	0.46
23:BB:1313:U:C2'	23:BB:1313:U:O2	2.64	0.46
23:BB:1408:G:H2'	23:BB:1409:U:H6	1.80	0.46
23:BB:1553:A:O2'	23:BB:1554:U:H2'	2.16	0.46
23:BB:1704:C:O2'	23:BB:1705:A:H5'	2.16	0.46
23:BB:192:C:C2'	23:BB:193:U:H5'	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1979:U:C2'	23:BB:1980:G:H5'	2.46	0.46
23:BB:1999:C:O2'	23:BB:2000:C:H5'	2.15	0.46
23:BB:2014:A:P	40:BS:95:ARG:HH21	2.39	0.46
23:BB:2307:G:O2'	23:BB:2311:A:N6	2.49	0.46
23:BB:2578:G:H1'	26:BD:145:SER:CB	2.46	0.46
23:BB:2600:A:H62	25:BC:235:GLU:CB	2.29	0.46
23:BB:441:U:H2'	23:BB:442:G:C8	2.49	0.46
23:BB:523:C:O2'	23:BB:524:G:H5'	2.16	0.46
23:BB:593:U:H2'	23:BB:594:U:H6	1.81	0.46
23:BB:85:G:H5'	42:BU:28:LEU:CA	2.22	0.46
23:BB:89:A:H2'	23:BB:90:U:C6	2.51	0.46
23:BB:935:C:H2'	23:BB:936:A:H8	1.81	0.46
25:BC:66:PHE:O	25:BC:68:ARG:HG3	2.16	0.46
25:BC:77:VAL:HG21	25:BC:109:LEU:CG	2.43	0.46
26:BD:15:PHE:HB3	26:BD:18:ASP:CB	2.45	0.46
26:BD:106:LYS:O	26:BD:206:ALA:HB2	2.16	0.46
28:BF:131:VAL:HG12	28:BF:132:ARG:N	2.30	0.46
28:BF:109:ARG:CB	28:BF:136:ILE:HG13	2.46	0.46
28:BF:56:LEU:HD23	28:BF:59:ILE:HG22	1.98	0.46
29:BG:34:ARG:HG3	29:BG:35:THR:N	2.30	0.46
30:BH:27:ARG:HB2	30:BH:27:ARG:HH11	1.81	0.46
52:BI:52:LEU:HD21	52:BI:81:LYS:NZ	2.31	0.46
32:BK:58:LEU:HA	32:BK:89:ASN:HD22	1.80	0.46
33:BL:103:ILE:HD13	33:BL:104:GLN:H	1.81	0.46
33:BL:56:PRO:C	33:BL:60:ARG:HB2	2.35	0.46
33:BL:94:THR:HG23	33:BL:95:LEU:H	1.80	0.46
36:BO:106:LEU:O	36:BO:106:LEU:HD13	2.16	0.46
23:BB:2334:U:OP2	36:BO:7:ARG:NE	2.49	0.46
37:BP:109:ILE:HG22	37:BP:109:ILE:O	2.16	0.46
26:BD:17:GLU:HA	37:BP:80:VAL:CG2	2.46	0.46
39:BR:48:LYS:C	39:BR:49:ILE:HD13	2.36	0.46
40:BS:4:ILE:HD12	40:BS:106:VAL:HG22	1.98	0.46
40:BS:28:LYS:HB3	40:BS:71:VAL:HB	1.98	0.46
40:BS:8:ARG:O	40:BS:9:HIS:HB2	2.16	0.46
42:BU:66:VAL:O	42:BU:68:ASN:N	2.48	0.46
42:BU:90:LYS:HD2	42:BU:91:LYS:H	1.81	0.46
24:BV:70:ILE:HG12	24:BV:71:LYS:N	2.27	0.46
43:BW:43:LYS:N	43:BW:43:LYS:HD2	2.31	0.46
46:BZ:7:PRO:O	46:BZ:8:LYS:HB3	2.16	0.46
1:CA:1049:U:H2'	1:CA:1049:U:O2	2.14	0.46
1:CA:1216:A:P	13:CN:2:LYS:HE2	2.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1279:G:H22	9:CJ:45:ARG:HD2	1.80	0.46
1:CA:132:C:O2'	1:CA:133:U:H5'	2.15	0.46
1:CA:22:G:O2'	1:CA:23:C:H5'	2.16	0.46
1:CA:411:A:C5	1:CA:413:G:H1'	2.51	0.46
1:CA:652:U:H1'	1:CA:653:U:C5	2.50	0.46
1:CA:861:G:O2'	1:CA:862:C:H5'	2.15	0.46
1:CA:932:C:H2'	1:CA:933:G:H8	1.80	0.46
1:CA:956:U:H2'	1:CA:957:U:C6	2.50	0.46
20:CB:195:VAL:HG12	20:CB:197:PHE:H	1.81	0.46
20:CB:60:ALA:HA	20:CB:64:GLY:HA3	1.98	0.46
2:CC:143:LEU:N	2:CC:143:LEU:HD13	2.31	0.46
4:CE:20:VAL:CG1	4:CE:31:SER:HB2	2.45	0.46
8:CI:117:LEU:HD23	8:CI:123:ARG:HD3	1.97	0.46
10:CK:17:ASP:CG	10:CK:36:ARG:HH11	2.19	0.46
12:CM:52:ILE:O	12:CM:55:LEU:HB2	2.16	0.46
14:CO:10:ILE:HD11	14:CO:29:ALA:CB	2.46	0.46
14:CO:81:ILE:O	14:CO:85:GLY:N	2.48	0.46
18:CS:32:THR:HG22	18:CS:33:TRP:N	2.21	0.46
48:D1:35:LEU:HD23	48:D1:35:LEU:N	2.31	0.46
50:D3:56:LEU:HD12	50:D3:56:LEU:C	2.36	0.46
51:D4:15:LYS:C	51:D4:17:VAL:H	2.20	0.46
23:DB:1597:A:H5''	23:DB:1598:A:H5'	1.96	0.46
23:DB:1665:A:O2'	23:DB:1666:G:H5'	2.15	0.46
23:DB:2065:C:H5'	23:DB:2251:G:H21	1.81	0.46
23:DB:2393:U:H5	50:D3:30:HIS:HE1	1.63	0.46
23:DB:2393:U:O2'	23:DB:2394:C:H5'	2.15	0.46
23:DB:2752:C:H2'	23:DB:2753:A:O4'	2.15	0.46
23:DB:2776:A:H4'	23:DB:2777:G:C5'	2.46	0.46
23:DB:2893:A:H4'	23:DB:2894:G:H5'	1.96	0.46
23:DB:494:G:O2'	23:DB:495:G:H5'	2.16	0.46
23:DB:741:U:H2'	23:DB:742:A:H8	1.79	0.46
31:DJ:28:LEU:HD22	31:DJ:28:LEU:O	2.16	0.46
31:DJ:57:LEU:HG	31:DJ:128:ASN:HA	1.98	0.46
33:DL:108:ALA:HB3	33:DL:125:LEU:HB2	1.98	0.46
35:DN:101:GLY:O	35:DN:102:PHE:HB3	2.16	0.46
35:DN:33:ILE:HG12	35:DN:114:GLU:HB3	1.98	0.46
37:DP:111:GLU:C	37:DP:113:LEU:H	2.18	0.46
38:DQ:40:LYS:O	38:DQ:44:TYR:HB3	2.16	0.46
40:DS:3:THR:HG21	40:DS:58:ALA:HB2	1.97	0.46
40:DS:5:ALA:HB3	40:DS:54:ALA:HB2	1.98	0.46
41:DT:47:VAL:HG22	41:DT:53:VAL:CG2	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DU:48:VAL:O	42:DU:48:VAL:HG12	2.16	0.46
46:DZ:64:PHE:HD2	46:DZ:64:PHE:O	1.98	0.46
1:AA:1098:C:C2'	1:AA:1099:G:H5'	2.46	0.45
1:AA:109:A:H4'	1:AA:110:C:OP2	2.16	0.45
1:AA:1149:C:H2'	1:AA:1150:A:H8	1.80	0.45
1:AA:1499:A:O2'	1:AA:1500:A:H5'	2.16	0.45
1:AA:525:C:H2'	1:AA:526:C:C6	2.51	0.45
1:AA:861:G:O2'	1:AA:862:C:H5'	2.15	0.45
1:AA:955:U:O2'	1:AA:956:U:H5'	2.16	0.45
20:AB:113:LEU:C	20:AB:113:LEU:HD23	2.37	0.45
20:AB:26:MET:HG2	20:AB:192:PRO:HG3	1.98	0.45
2:AC:102:ILE:C	2:AC:102:ILE:HD12	2.36	0.45
3:AD:160:LEU:H	3:AD:160:LEU:CD1	2.22	0.45
3:AD:169:TRP:CB	3:AD:183:ARG:HH21	2.30	0.45
5:AF:100:SER:HA	17:AR:23:LYS:HZ2	1.81	0.45
11:AL:90:PRO:C	11:AL:92:VAL:H	2.18	0.45
14:AO:70:LYS:HZ2	14:AO:74:VAL:HA	1.81	0.45
21:AU:32:ARG:HB3	21:AU:33:ARG:H	1.59	0.45
48:B1:12:SER:HA	48:B1:50:GLU:CB	2.41	0.45
50:B3:35:LYS:HD3	50:B3:35:LYS:N	2.31	0.45
23:BB:1064:C:O2'	23:BB:1065:U:H5'	2.16	0.45
23:BB:1123:C:H2'	23:BB:1124:G:C8	2.48	0.45
23:BB:1940:U:H1'	23:BB:1941:C:H5	1.81	0.45
23:BB:2721:A:H2'	23:BB:2722:G:C8	2.51	0.45
23:BB:359:G:C2'	23:BB:360:U:H5'	2.46	0.45
23:BB:532:A:N3	23:BB:532:A:H2'	2.30	0.45
23:BB:570:G:O2'	23:BB:571:U:H5'	2.16	0.45
23:BB:630:G:H4'	23:BB:640:C:O2'	2.16	0.45
23:BB:917:A:H2'	23:BB:918:A:O4'	2.15	0.45
23:BB:963:U:H2'	23:BB:964:C:C6	2.51	0.45
26:BD:50:VAL:HB	26:BD:75:ALA:HB3	1.97	0.45
27:BE:144:GLU:C	27:BE:146:VAL:N	2.70	0.45
28:BF:32:LYS:HB3	28:BF:91:ARG:HB3	1.98	0.45
29:BG:57:TYR:CZ	29:BG:59:ASP:HB2	2.52	0.45
23:BB:2748:A:H1'	29:BG:66:THR:OG1	2.16	0.45
29:BG:71:LEU:HG	29:BG:74:MET:SD	2.56	0.45
31:BJ:43:GLU:O	31:BJ:44:TYR:C	2.54	0.45
33:BL:77:ILE:HB	33:BL:110:VAL:HG13	1.98	0.45
33:BL:125:LEU:O	33:BL:126:ARG:HG3	2.17	0.45
34:BM:124:LEU:H	34:BM:124:LEU:HD23	1.81	0.45
36:BO:94:ARG:O	36:BO:95:SER:C	2.54	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:88:ARG:HA	37:BP:88:ARG:NE	2.30	0.45
39:BR:42:ALA:O	39:BR:45:GLU:HG3	2.16	0.45
40:BS:8:ARG:NH2	40:BS:102:HIS:NE2	2.64	0.45
23:BB:1157:G:H1'	45:BY:10:ARG:HH22	1.81	0.45
46:BZ:63:ARG:HA	46:BZ:63:ARG:NE	2.31	0.45
1:CA:1237:C:H3'	1:CA:1336:C:H41	1.81	0.45
1:CA:1250:A:H5''	8:CI:69:GLY:H	1.81	0.45
1:CA:202:G:N2	1:CA:465:A:H61	2.14	0.45
1:CA:357:G:OP1	1:CA:367:U:H6	1.99	0.45
1:CA:598:U:H2'	1:CA:599:C:C6	2.51	0.45
1:CA:767:A:H2'	1:CA:768:A:H8	1.81	0.45
1:CA:834:U:H2'	1:CA:835:U:C6	2.50	0.45
1:CA:924:C:H2'	1:CA:925:G:H8	1.82	0.45
20:CB:67:LEU:O	20:CB:160:LEU:HD12	2.16	0.45
2:CC:145:ALA:HA	2:CC:203:LYS:HA	1.98	0.45
3:CD:127:ARG:H	3:CD:127:ARG:CD	2.28	0.45
5:CF:67:PRO:O	5:CF:70:VAL:HG22	2.15	0.45
8:CI:94:ARG:O	8:CI:98:ARG:HG3	2.15	0.45
9:CJ:17:LEU:C	9:CJ:17:LEU:HD13	2.35	0.45
14:CO:36:ASN:HA	14:CO:39:GLN:HG3	1.98	0.45
16:CQ:56:ASP:N	16:CQ:81:ALA:HB2	2.31	0.45
1:CA:1014:A:H5''	18:CS:13:HIS:CB	2.46	0.45
49:D2:46:LYS:N	49:D2:46:LYS:HE3	2.31	0.45
23:DB:1033:U:C4	51:D4:16:ILE:HD11	2.52	0.45
23:DB:1084:A:H1'	23:DB:1106:G:H5'	1.98	0.45
23:DB:1678:A:O2'	23:DB:1679:A:H5'	2.16	0.45
23:DB:1747:U:H2'	23:DB:1748:C:H6	1.80	0.45
23:DB:786:C:H5''	23:DB:1780:A:N7	2.31	0.45
23:DB:1816:C:C5	25:DC:62:ARG:HD2	2.51	0.45
23:DB:2087:G:H2'	23:DB:2088:A:C8	2.51	0.45
23:DB:2182:U:H2'	23:DB:2183:A:C8	2.52	0.45
23:DB:2379:G:H2'	23:DB:2380:C:C6	2.51	0.45
23:DB:2477:U:H4'	23:DB:2479:U:O4	2.16	0.45
23:DB:2746:U:O2'	23:DB:2747:G:H5'	2.16	0.45
23:DB:299:A:N6	23:DB:322:A:H1'	2.31	0.45
23:DB:346:A:H2'	23:DB:347:A:O4'	2.15	0.45
23:DB:383:C:H41	23:DB:385:C:H2'	1.81	0.45
23:DB:448:U:C6	27:DE:79:ARG:HG3	2.51	0.45
25:DC:142:ASN:HA	25:DC:153:LEU:CD2	2.46	0.45
23:DB:1821:A:C5'	25:DC:155:ARG:HH21	2.29	0.45
26:DD:138:LEU:HD12	26:DD:142:VAL:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:DE:10:SER:C	27:DE:12:LEU:H	2.19	0.45
27:DE:55:SER:N	27:DE:74:LYS:HE3	2.30	0.45
32:DK:12:ASP:HA	32:DK:99:ILE:HA	1.97	0.45
33:DL:18:ARG:NH2	33:DL:21:ARG:HD3	2.31	0.45
23:DB:873:C:H4'	34:DM:64:TRP:CZ3	2.51	0.45
34:DM:85:GLY:O	34:DM:86:LYS:O	2.34	0.45
37:DP:52:ARG:HG3	37:DP:54:LEU:HD12	1.98	0.45
41:DT:20:ALA:O	41:DT:21:SER:HB3	2.17	0.45
42:DU:48:VAL:C	42:DU:50:ALA:N	2.65	0.45
43:DW:77:LYS:C	43:DW:79:ILE:N	2.68	0.45
45:DY:25:GLY:C	45:DY:26:LEU:HD23	2.36	0.45
45:DY:37:ARG:NE	45:DY:37:ARG:HA	2.31	0.45
1:AA:1430:A:H2'	1:AA:1431:A:O4'	2.17	0.45
1:AA:1473:G:H2'	1:AA:1474:U:O4'	2.16	0.45
1:AA:22:G:O2'	1:AA:23:C:H5'	2.17	0.45
1:AA:336:A:O2'	1:AA:337:G:H5'	2.15	0.45
1:AA:411:A:C5	1:AA:413:G:H1'	2.51	0.45
1:AA:767:A:H2'	1:AA:768:A:H8	1.81	0.45
1:AA:880:C:H2'	1:AA:881:G:H8	1.81	0.45
2:AC:70:ALA:HB2	2:AC:108:PRO:HB3	1.99	0.45
3:AD:56:GLU:HG2	3:AD:198:LEU:HD12	1.97	0.45
3:AD:64:TYR:H	3:AD:64:TYR:HD1	1.64	0.45
5:AF:7:VAL:O	5:AF:7:VAL:HG13	2.16	0.45
6:AG:11:ILE:HD12	6:AG:11:ILE:H	1.80	0.45
7:AH:104:SER:O	7:AH:122:GLY:HA3	2.16	0.45
7:AH:75:GLN:O	7:AH:126:CYS:HB2	2.17	0.45
10:AK:22:ILE:HD12	10:AK:85:VAL:HG22	1.98	0.45
11:AL:31:GLY:HA3	11:AL:54:VAL:CG1	2.46	0.45
15:AP:45:GLU:HG2	15:AP:45:GLU:O	2.16	0.45
18:AS:11:ASP:HB2	18:AS:14:LEU:CD2	2.47	0.45
23:BB:2056:G:O2'	47:B0:2:VAL:HG11	2.17	0.45
23:BB:125:A:H5'	49:B2:19:ARG:HE	1.81	0.45
23:BB:1112:G:O2'	29:BG:2:ARG:HG2	2.17	0.45
23:BB:1133:A:H4'	23:BB:1134:A:C5'	2.46	0.45
23:BB:1164:C:H2'	23:BB:1165:A:H8	1.81	0.45
23:BB:1184:U:H2'	23:BB:1185:G:H8	1.81	0.45
23:BB:1541:C:H2'	23:BB:1542:U:O4'	2.16	0.45
23:BB:1597:A:H5''	23:BB:1598:A:H5'	1.97	0.45
23:BB:1760:C:H2'	23:BB:1761:C:O4'	2.16	0.45
23:BB:1790:C:H2'	23:BB:1791:A:C8	2.51	0.45
23:BB:1803:A:H2	23:BB:1823:G:C1'	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1915:U:H3'	23:BB:1916:A:C8	2.50	0.45
23:BB:528:A:N1	23:BB:2043:C:H5'	2.31	0.45
23:BB:764:A:N1	23:BB:1789:A:O2'	2.47	0.45
26:BD:146:ILE:HG21	26:BD:159:LYS:NZ	2.30	0.45
26:BD:184:ARG:HG3	26:BD:184:ARG:O	2.16	0.45
26:BD:29:VAL:HA	26:BD:185:ASN:CG	2.37	0.45
29:BG:104:LEU:HD12	29:BG:104:LEU:N	2.31	0.45
29:BG:75:VAL:HG23	29:BG:76:ILE:N	2.31	0.45
52:BI:121:ILE:O	52:BI:124:MET:HG2	2.16	0.45
31:BJ:31:GLU:O	31:BJ:35:ARG:HB2	2.15	0.45
33:BL:77:ILE:HG22	33:BL:78:ARG:N	2.30	0.45
23:BB:2276:G:OP2	34:BM:83:GLY:HA3	2.16	0.45
42:BU:3:LYS:HD2	42:BU:81:ARG:HH22	1.81	0.45
1:CA:744:C:H2'	1:CA:745:G:C8	2.50	0.45
20:CB:95:TRP:CZ2	20:CB:100:LEU:HD13	2.50	0.45
2:CC:38:VAL:O	2:CC:42:LEU:HD23	2.17	0.45
2:CC:76:ILE:HA	2:CC:83:VAL:CG2	2.46	0.45
2:CC:94:ALA:O	2:CC:96:VAL:N	2.50	0.45
3:CD:47:LEU:HD11	3:CD:52:VAL:N	2.31	0.45
3:CD:54:LEU:O	3:CD:58:GLN:HB2	2.17	0.45
4:CE:156:ARG:NH2	7:CH:100:ILE:HG23	2.32	0.45
6:CG:41:ILE:HD13	6:CG:115:MET:HG3	1.97	0.45
7:CH:45:ILE:CD1	7:CH:60:LEU:HD11	2.46	0.45
9:CJ:17:LEU:HD22	9:CJ:17:LEU:O	2.16	0.45
9:CJ:8:ILE:HD11	9:CJ:74:VAL:HB	1.98	0.45
14:CO:10:ILE:HG21	14:CO:30:LEU:CD2	2.45	0.45
16:CQ:10:ARG:CZ	16:CQ:55:GLY:N	2.79	0.45
1:CA:720:C:O4'	17:CR:38:ILE:HG12	2.16	0.45
22:DA:55:U:H2'	22:DA:56:G:C8	2.51	0.45
23:DB:1353:A:H2'	23:DB:1354:A:C8	2.50	0.45
23:DB:1683:U:H2'	23:DB:1684:G:C8	2.51	0.45
23:DB:1704:C:O2'	23:DB:1705:A:H5'	2.16	0.45
23:DB:1910:G:O2'	23:DB:1911:U:H5'	2.16	0.45
23:DB:1979:U:C2'	23:DB:1980:G:H5'	2.45	0.45
23:DB:2038:G:H2'	23:DB:2039:U:O4'	2.17	0.45
23:DB:257:C:H2'	23:DB:258:G:O4'	2.17	0.45
23:DB:264:C:H2'	23:DB:265:A:H5''	1.98	0.45
23:DB:315:G:O2'	23:DB:316:C:H5'	2.16	0.45
23:DB:439:A:H2'	23:DB:440:C:O4'	2.16	0.45
23:DB:630:G:H4'	23:DB:640:C:O2'	2.16	0.45
23:DB:692:C:H2'	23:DB:693:A:H8	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:171:VAL:O	25:DC:172:THR:HG23	2.17	0.45
25:DC:178:GLY:C	25:DC:179:GLU:HG2	2.35	0.45
25:DC:235:GLU:CG	25:DC:237:ARG:HD3	2.46	0.45
29:DG:55:ASP:CG	29:DG:56:GLY:N	2.69	0.45
30:DH:109:GLU:O	30:DH:109:GLU:HG3	2.16	0.45
30:DH:126:GLY:H	30:DH:146:VAL:HB	1.82	0.45
32:DK:78:ARG:H	37:DP:72:VAL:CG2	2.27	0.45
32:DK:87:LEU:HD22	32:DK:92:GLU:O	2.17	0.45
33:DL:77:ILE:HD13	33:DL:110:VAL:CA	2.44	0.45
35:DN:48:VAL:O	35:DN:51:LEU:HB2	2.15	0.45
36:DO:30:ARG:CG	36:DO:31:THR:H	2.28	0.45
36:DO:27:VAL:HG22	36:DO:38:GLN:O	2.17	0.45
39:DR:4:VAL:O	39:DR:41:ILE:HG21	2.16	0.45
38:DQ:88:GLU:OE2	39:DR:53:PHE:HB2	2.16	0.45
40:DS:87:PRO:O	40:DS:88:ARG:CB	2.64	0.45
41:DT:38:ALA:HB1	41:DT:43:ILE:CD1	2.42	0.45
42:DU:43:LYS:NZ	42:DU:43:LYS:HB2	2.30	0.45
46:DZ:37:CYS:SG	46:DZ:39:LYS:HB2	2.57	0.45
1:AA:1206:G:H4'	2:AC:192:TYR:HA	1.98	0.45
1:AA:1490:U:H5'	1:AA:1491:G:OP2	2.16	0.45
1:AA:167:A:H2'	1:AA:168:G:H8	1.81	0.45
1:AA:299:G:H2'	1:AA:300:A:C8	2.51	0.45
12:AM:102:LYS:NZ	12:AM:102:LYS:HB2	2.31	0.45
12:AM:47:LEU:HB3	12:AM:51:GLN:HB2	1.97	0.45
17:AR:62:ARG:HD3	17:AR:69:TYR:HA	1.99	0.45
18:AS:43:MET:O	18:AS:46:LEU:HB2	2.16	0.45
50:B3:12:ARG:CZ	50:B3:12:ARG:HB2	2.41	0.45
50:B3:16:THR:HG21	50:B3:22:LYS:NZ	2.32	0.45
51:B4:7:VAL:HG12	51:B4:8:LYS:N	2.25	0.45
23:BB:1289:C:H2'	23:BB:1290:C:C6	2.51	0.45
23:BB:1720:U:C2'	23:BB:1721:G:H5'	2.46	0.45
23:BB:2261:C:O2'	23:BB:2262:U:H5'	2.15	0.45
23:BB:2800:A:C4	23:BB:2801:G:H1'	2.51	0.45
23:BB:28:A:H1'	23:BB:513:A:C2	2.52	0.45
23:BB:492:A:H2'	23:BB:493:G:O4'	2.16	0.45
23:BB:997:G:C5'	38:BQ:91:ARG:HG3	2.47	0.45
23:BB:1818:U:H2'	25:BC:152:GLN:O	2.17	0.45
25:BC:140:VAL:HA	25:BC:191:LEU:HA	1.98	0.45
25:BC:198:GLU:C	25:BC:200:MET:H	2.18	0.45
25:BC:66:PHE:O	25:BC:67:LYS:C	2.55	0.45
25:BC:68:ARG:HD2	25:BC:127:ASN:ND2	2.27	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2512:C:O2'	26:BD:146:ILE:HG12	2.15	0.45
27:BE:6:LYS:HB2	27:BE:12:LEU:CD1	2.46	0.45
27:BE:146:VAL:HG23	27:BE:148:ILE:CG2	2.46	0.45
28:BF:15:LEU:HD13	28:BF:28:PRO:CG	2.45	0.45
28:BF:33:ILE:HG22	28:BF:155:ILE:HA	1.98	0.45
30:BH:1:MET:C	30:BH:21:VAL:HG12	2.37	0.45
31:BJ:25:LEU:HD13	31:BJ:25:LEU:H	1.81	0.45
31:BJ:50:THR:OG1	31:BJ:51:GLY:N	2.49	0.45
31:BJ:25:LEU:CD1	31:BJ:62:VAL:HA	2.41	0.45
34:BM:8:LYS:NZ	34:BM:70:ASP:H	2.10	0.45
35:BN:86:ARG:NH2	35:BN:116:VAL:HG11	2.31	0.45
35:BN:9:GLN:NE2	35:BN:17:ARG:NH2	2.65	0.45
36:BO:31:THR:HG21	36:BO:36:TYR:HE1	1.82	0.45
37:BP:60:VAL:HB	37:BP:61:ARG:H	1.46	0.45
40:BS:102:HIS:O	40:BS:103:ILE:HB	2.15	0.45
40:BS:69:LEU:HA	40:BS:108:SER:O	2.16	0.45
41:BT:57:VAL:O	41:BT:85:VAL:HA	2.16	0.45
24:BV:44:HIS:O	24:BV:48:MET:HB3	2.16	0.45
43:BW:58:LEU:HA	43:BW:79:ILE:HG22	1.98	0.45
45:BY:29:ARG:O	45:BY:30:ARG:HB2	2.17	0.45
1:CA:1000:A:O2'	1:CA:1001:C:H5'	2.16	0.45
1:CA:1004:A:N7	1:CA:1025:U:O2'	2.48	0.45
1:CA:1013:G:N2	1:CA:1015:G:H3'	2.32	0.45
1:CA:1054:C:O2'	1:CA:1055:A:H5''	2.17	0.45
1:CA:1124:G:H5'	9:CJ:37:ARG:HH21	1.81	0.45
1:CA:1187:G:H2'	1:CA:1188:A:O4'	2.16	0.45
1:CA:1456:A:H2'	1:CA:1457:G:H8	1.81	0.45
1:CA:167:A:H2'	1:CA:168:G:H8	1.81	0.45
1:CA:299:G:H2'	1:CA:300:A:C8	2.51	0.45
1:CA:393:A:H5'	1:CA:483:C:O2'	2.15	0.45
1:CA:977:A:H8	1:CA:1223:C:N3	2.15	0.45
20:CB:209:VAL:HG23	20:CB:210:THR:N	2.31	0.45
20:CB:60:ALA:HB2	20:CB:66:ILE:HD11	1.98	0.45
2:CC:99:GLN:O	2:CC:100:ILE:HB	2.16	0.45
3:CD:30:LYS:HD3	3:CD:30:LYS:N	2.31	0.45
1:CA:7:A:H2	4:CE:125:LYS:HE2	1.82	0.45
4:CE:131:ASN:C	4:CE:135:VAL:HG23	2.35	0.45
8:CI:56:MET:HA	8:CI:59:LYS:HB2	1.98	0.45
9:CJ:36:VAL:HG22	9:CJ:76:ILE:HD13	1.98	0.45
10:CK:126:ARG:HG2	10:CK:126:ARG:HH11	1.80	0.45
12:CM:79:LEU:HB2	12:CM:84:CYS:SG	2.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1308:U:OP2	12:CM:97:ARG:HD3	2.16	0.45
15:CP:32:PHE:CD1	15:CP:32:PHE:C	2.90	0.45
16:CQ:74:LEU:HD22	16:CQ:75:VAL:H	1.82	0.45
17:CR:61:ALA:HB3	17:CR:67:LEU:HD12	1.98	0.45
19:CT:19:HIS:CE1	19:CT:23:ARG:HG3	2.51	0.45
22:DA:12:C:H4'	22:DA:15:A:N6	2.32	0.45
22:DA:63:C:H2'	22:DA:64:G:C8	2.51	0.45
23:DB:1047:G:H1'	23:DB:1110:G:N2	2.32	0.45
23:DB:1289:C:H2'	23:DB:1290:C:C6	2.51	0.45
23:DB:1313:U:O2	23:DB:1313:U:C2'	2.64	0.45
23:DB:1322:A:H2'	23:DB:1323:C:H5'	1.98	0.45
23:DB:137:U:O2'	23:DB:138:U:H5'	2.16	0.45
23:DB:1601:G:O2'	23:DB:1602:U:H5'	2.16	0.45
23:DB:1668:A:N3	23:DB:1670:C:C4	2.84	0.45
23:DB:1708:C:H2'	23:DB:1709:U:H6	1.82	0.45
23:DB:2103:C:H3'	23:DB:2104:C:C2	2.51	0.45
23:DB:2147:A:C2'	23:DB:2147:A:N3	2.78	0.45
23:DB:2207:C:H2'	23:DB:2208:C:H6	1.82	0.45
23:DB:2465:C:O2'	23:DB:2466:C:H5'	2.17	0.45
23:DB:265:A:O2'	23:DB:266:G:H4'	2.17	0.45
23:DB:2772:C:H2'	23:DB:2773:C:H6	1.81	0.45
23:DB:2830:C:O2'	23:DB:2831:G:H5'	2.17	0.45
23:DB:345:A:N3	23:DB:346:A:N1	2.64	0.45
23:DB:522:A:H2'	23:DB:523:C:H6	1.80	0.45
23:DB:709:U:H2'	23:DB:710:U:C6	2.51	0.45
23:DB:780:G:OP1	25:DC:216:ARG:NH2	2.49	0.45
23:DB:992:C:H4'	38:DQ:46:TYR:HH	1.81	0.45
25:DC:19:VAL:CG1	25:DC:20:ASN:N	2.80	0.45
27:DE:4:VAL:HA	27:DE:14:VAL:CG1	2.46	0.45
28:DF:108:PRO:O	28:DF:109:ARG:HB2	2.17	0.45
28:DF:70:ARG:HA	28:DF:80:GLN:NE2	2.20	0.45
29:DG:131:VAL:O	29:DG:131:VAL:HG13	2.15	0.45
29:DG:171:LYS:HB2	29:DG:172:GLU:H	1.61	0.45
30:DH:92:GLY:O	30:DH:93:SER:HB3	2.16	0.45
52:DI:96:LYS:HD3	52:DI:138:VAL:HG21	1.96	0.45
23:DB:1099:G:C4'	52:DI:4:VAL:HB	2.43	0.45
52:DI:9:LYS:HG2	52:DI:57:VAL:HG13	1.98	0.45
35:DN:16:HIS:O	35:DN:20:MET:N	2.41	0.45
36:DO:25:ARG:HG2	36:DO:94:ARG:NH2	2.27	0.45
38:DQ:52:ARG:HG2	38:DQ:53:LYS:CE	2.46	0.45
40:DS:46:LEU:CA	40:DS:49:LYS:HB2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DS:85:ILE:HG12	40:DS:94:ASP:O	2.16	0.45
41:DT:68:LYS:HZ3	41:DT:68:LYS:HB2	1.79	0.45
41:DT:30:ILE:O	41:DT:85:VAL:HG22	2.17	0.45
24:DV:93:ARG:HG2	24:DV:94:ALA:H	1.82	0.45
45:DY:6:ILE:O	45:DY:6:ILE:HG13	2.16	0.45
46:DZ:25:ARG:CG	46:DZ:26:SER:H	2.22	0.45
1:AA:1074:G:H5'	20:AB:104:LYS:HZ1	1.79	0.45
1:AA:1258:G:C4	1:AA:1278:G:N2	2.84	0.45
1:AA:708:C:H2'	1:AA:709:U:H6	1.79	0.45
1:AA:744:C:H2'	1:AA:745:G:C8	2.50	0.45
1:AA:836:G:H2'	1:AA:837:U:C6	2.52	0.45
1:AA:962:C:H2'	1:AA:963:G:O4'	2.17	0.45
20:AB:72:LYS:NZ	20:AB:167:HIS:HB2	2.32	0.45
20:AB:204:ASP:CG	20:AB:205:ALA:N	2.70	0.45
20:AB:96:LEU:HD21	20:AB:146:SER:HB2	1.98	0.45
2:AC:186:SER:O	2:AC:197:VAL:HG12	2.16	0.45
3:AD:163:GLN:N	3:AD:163:GLN:OE1	2.50	0.45
4:AE:81:GLN:NE2	4:AE:147:ASN:N	2.65	0.45
7:AH:48:PHE:HB2	7:AH:60:LEU:HD12	1.99	0.45
10:AK:112:VAL:HA	17:AR:72:ARG:HD3	1.99	0.45
14:AO:78:THR:HA	14:AO:81:ILE:CD1	2.46	0.45
15:AP:29:ASN:N	15:AP:29:ASN:ND2	2.63	0.45
15:AP:71:VAL:HA	15:AP:74:LEU:CG	2.45	0.45
21:AU:43:GLU:HA	21:AU:43:GLU:OE1	2.15	0.45
47:B0:54:ILE:O	47:B0:54:ILE:HG13	2.17	0.45
48:B1:41:VAL:HB	48:B1:43:ARG:NH2	2.31	0.45
50:B3:26:ALA:HB3	50:B3:29:ARG:CZ	2.46	0.45
22:BA:29:A:H3'	22:BA:30:C:C6	2.49	0.45
22:BA:64:G:O2'	22:BA:65:U:H5'	2.17	0.45
23:BB:1665:A:O2'	23:BB:1666:G:H5'	2.15	0.45
23:BB:2038:G:H2'	23:BB:2039:U:O4'	2.17	0.45
23:BB:2038:G:H2'	23:BB:2039:U:H6	1.82	0.45
23:BB:189:G:H2'	23:BB:205:G:H22	1.81	0.45
23:BB:2301:C:O2'	23:BB:2302:U:H5'	2.17	0.45
23:BB:2363:G:O2'	23:BB:2364:C:H5'	2.17	0.45
23:BB:2556:C:O2'	23:BB:2557:G:H5'	2.17	0.45
23:BB:327:G:O2'	23:BB:328:U:H5'	2.16	0.45
23:BB:407:G:O2'	23:BB:408:G:H5'	2.16	0.45
23:BB:48:G:H4'	23:BB:52:A:O4'	2.16	0.45
23:BB:760:G:H4'	23:BB:1776:G:OP1	2.16	0.45
25:BC:136:VAL:HA	25:BC:165:ALA:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:161:VAL:H	25:BC:193:GLU:CD	2.20	0.45
25:BC:260:LYS:C	25:BC:262:THR:N	2.70	0.45
26:BD:115:GLY:HA2	26:BD:165:MET:CG	2.35	0.45
26:BD:12:THR:HB	26:BD:21:SER:O	2.15	0.45
27:BE:194:LYS:HE3	27:BE:201:ALA:HB2	1.98	0.45
29:BG:51:PHE:H	29:BG:68:ARG:HD3	1.80	0.45
29:BG:25:ILE:HD11	29:BG:75:VAL:HG12	1.97	0.45
30:BH:50:ARG:HG2	30:BH:50:ARG:O	2.15	0.45
23:BB:1060:U:C4	52:BI:131:THR:HG22	2.51	0.45
31:BJ:120:ARG:O	31:BJ:122:LEU:N	2.42	0.45
23:BB:6:A:H5''	31:BJ:133:ALA:O	2.17	0.45
31:BJ:52:ASP:CG	31:BJ:53:TYR:H	2.19	0.45
31:BJ:59:ALA:C	31:BJ:61:LYS:H	2.20	0.45
33:BL:82:LEU:HG	33:BL:122:VAL:HG12	1.99	0.45
35:BN:99:LYS:HZ2	35:BN:99:LYS:HB3	1.77	0.45
36:BO:97:PHE:O	36:BO:98:GLN:HG2	2.16	0.45
38:BQ:12:ARG:HH11	38:BQ:12:ARG:HG2	1.82	0.45
39:BR:14:VAL:HG12	39:BR:15:SER:N	2.32	0.45
39:BR:20:VAL:HG22	39:BR:97:LYS:HZ1	1.81	0.45
39:BR:53:PHE:HD2	39:BR:55:ASP:H	1.59	0.45
40:BS:88:ARG:HB3	40:BS:89:ALA:H	1.57	0.45
46:BZ:19:GLY:O	46:BZ:20:ASN:HB2	2.16	0.45
1:CA:29:U:H5'	1:CA:296:U:OP1	2.17	0.45
1:CA:328:C:H4'	1:CA:329:A:H5''	1.98	0.45
1:CA:525:C:H2'	1:CA:526:C:C6	2.51	0.45
1:CA:812:G:C2'	1:CA:812:G:N3	2.77	0.45
20:CB:132:GLU:HA	20:CB:135:MET:CE	2.47	0.45
3:CD:141:VAL:HG12	3:CD:180:THR:CA	2.45	0.45
10:CK:60:PHE:O	10:CK:64:VAL:HG13	2.15	0.45
11:CL:36:VAL:HG12	11:CL:52:CYS:HB2	1.96	0.45
1:CA:1309:G:C5'	12:CM:76:ILE:HG12	2.46	0.45
16:CQ:46:HIS:CG	16:CQ:47:ASP:H	2.33	0.45
5:CF:100:SER:N	17:CR:23:LYS:NZ	2.63	0.45
19:CT:34:VAL:CG1	19:CT:78:LEU:HD22	2.47	0.45
21:CU:23:GLU:HA	21:CU:27:VAL:HG21	1.98	0.45
21:CU:39:LYS:N	21:CU:40:PRO:HD2	2.21	0.45
23:DB:1259:G:O2'	23:DB:1260:A:H5'	2.17	0.45
23:DB:1408:G:H2'	23:DB:1409:U:H6	1.80	0.45
23:DB:1553:A:O2'	23:DB:1554:U:H2'	2.16	0.45
23:DB:2223:G:O2'	23:DB:2224:G:H5'	2.16	0.45
23:DB:2243:U:O2	23:DB:2434:A:C2	2.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2276:G:OP1	34:DM:82:MET:SD	2.75	0.45
23:DB:2529:G:H4'	29:DG:175:LYS:HD3	1.98	0.45
23:DB:2714:G:H2'	23:DB:2715:C:C6	2.51	0.45
23:DB:359:G:C2'	23:DB:360:U:H5'	2.46	0.45
23:DB:534:U:H5'	38:DQ:41:ALA:CA	2.43	0.45
23:DB:817:C:O2'	23:DB:839:U:H5''	2.17	0.45
25:DC:152:GLN:HB3	25:DC:153:LEU:H	1.62	0.45
25:DC:227:VAL:HG13	25:DC:228:ASP:OD1	2.17	0.45
27:DE:114:ARG:HG3	27:DE:114:ARG:NH1	2.32	0.45
27:DE:23:PHE:CG	27:DE:24:ASN:N	2.85	0.45
30:DH:4:ILE:HD11	30:DH:37:VAL:O	2.17	0.45
23:DB:1099:G:H5''	52:DI:2:LYS:HB2	1.98	0.45
31:DJ:85:LYS:NZ	31:DJ:86:GLN:O	2.49	0.45
33:DL:50:PHE:O	33:DL:52:GLY:N	2.49	0.45
33:DL:55:MET:SD	33:DL:56:PRO:HG3	2.56	0.45
35:DN:52:ILE:HD11	35:DN:83:LEU:HD21	1.99	0.45
35:DN:96:ARG:NH1	35:DN:98:LEU:HD21	2.31	0.45
36:DO:30:ARG:O	36:DO:31:THR:HB	2.16	0.45
39:DR:99:THR:HG23	39:DR:100:GLY:H	1.80	0.45
39:DR:76:LYS:HB3	39:DR:90:ARG:HG2	1.96	0.45
40:DS:22:ASP:C	40:DS:24:ILE:H	2.20	0.45
40:DS:81:SER:HB2	40:DS:99:ARG:H	1.81	0.45
43:DW:36:ILE:HG21	43:DW:68:PHE:HE1	1.79	0.45
1:AA:1014:A:OP2	18:AS:17:LYS:NZ	2.45	0.45
1:AA:1113:C:H2'	1:AA:1114:C:H6	1.82	0.45
1:AA:1197:A:P	1:AA:1197:A:H3'	2.57	0.45
1:AA:1410:A:H2'	1:AA:1411:C:H6	1.79	0.45
1:AA:29:U:O2'	1:AA:30:U:H5'	2.17	0.45
1:AA:972:C:H4'	9:AJ:59:LYS:HB3	1.99	0.45
20:AB:9:LEU:CD2	20:AB:11:ALA:H	2.17	0.45
2:AC:112:ALA:HB1	2:AC:199:VAL:HG23	1.99	0.45
3:AD:98:ASP:CG	3:AD:132:ALA:HB1	2.36	0.45
5:AF:1:MET:HB3	5:AF:65:GLU:O	2.16	0.45
6:AG:71:THR:HA	6:AG:90:VAL:HG22	1.99	0.45
8:AI:46:VAL:HB	8:AI:79:ARG:HD3	1.97	0.45
1:AA:718:A:H5'	10:AK:118:ASN:HB2	1.97	0.45
12:AM:21:ILE:HG22	12:AM:64:VAL:CG1	2.44	0.45
15:AP:7:ALA:HB1	15:AP:29:ASN:CB	2.47	0.45
18:AS:4:LEU:O	18:AS:6:LYS:N	2.49	0.45
48:B1:42:VAL:CG1	48:B1:43:ARG:HE	2.28	0.45
22:BA:18:G:H2'	22:BA:19:C:C6	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1125:G:H3'	23:BB:1126:A:H5''	1.98	0.45
23:BB:137:U:O5'	23:BB:137:U:H6	1.99	0.45
23:BB:1410:G:H2'	23:BB:1411:U:H6	1.77	0.45
23:BB:1450:G:N2	23:BB:1452:G:H1	1.95	0.45
23:BB:1708:C:H2'	23:BB:1709:U:H6	1.81	0.45
23:BB:1885:A:H3'	23:BB:1886:U:C6	2.51	0.45
23:BB:215:G:O4'	23:BB:216:A:H4'	2.17	0.45
23:BB:2187:U:H2'	23:BB:2187:U:O2	2.16	0.45
23:BB:2710:C:H2'	23:BB:2711:A:H8	1.80	0.45
23:BB:2722:G:H2'	23:BB:2723:C:H6	1.82	0.45
23:BB:2751:G:OP2	29:BG:3:VAL:HG13	2.16	0.45
23:BB:2780:G:C2	31:BJ:120:ARG:HG3	2.51	0.45
23:BB:364:C:O2'	23:BB:365:U:H5'	2.16	0.45
23:BB:539:G:H2'	23:BB:540:C:H6	1.81	0.45
23:BB:773:U:H4'	25:BC:47:ARG:CA	2.47	0.45
25:BC:163:ILE:HG22	25:BC:172:THR:HG21	1.98	0.45
26:BD:165:MET:HG3	26:BD:166:GLY:N	2.32	0.45
33:BL:120:VAL:CG1	33:BL:138:ALA:HB3	2.47	0.45
35:BN:32:GLU:O	35:BN:114:GLU:HA	2.17	0.45
39:BR:66:HIS:CA	39:BR:98:ILE:HD13	2.45	0.45
40:BS:29:VAL:C	40:BS:31:GLN:N	2.69	0.45
41:BT:25:GLU:CG	41:BT:26:LYS:N	2.79	0.45
41:BT:15:HIS:CB	41:BT:31:VAL:HG21	2.46	0.45
42:BU:3:LYS:HD2	42:BU:81:ARG:NH2	2.31	0.45
44:BX:16:THR:O	44:BX:20:ASN:HB2	2.17	0.45
44:BX:22:LEU:HA	44:BX:47:ARG:NH2	2.32	0.45
46:BZ:50:ASP:O	46:BZ:53:THR:HG22	2.16	0.45
1:CA:1060:U:C5'	9:CJ:53:ILE:HG12	2.46	0.45
1:CA:1065:U:H1'	1:CA:1066:C:OP2	2.17	0.45
1:CA:12:U:H4'	1:CA:526:C:H4'	1.98	0.45
1:CA:1425:U:H2'	1:CA:1426:G:H8	1.82	0.45
1:CA:613:C:H2'	1:CA:614:C:H6	1.82	0.45
1:CA:820:U:H4'	1:CA:821:G:OP2	2.16	0.45
1:CA:880:C:H2'	1:CA:881:G:H8	1.81	0.45
1:CA:950:U:H2'	1:CA:951:G:H8	1.82	0.45
1:CA:975:A:H2'	1:CA:976:G:OP2	2.17	0.45
20:CB:95:TRP:HZ2	20:CB:100:LEU:HD13	1.81	0.45
4:CE:125:LYS:HE3	4:CE:127:TYR:CE1	2.52	0.45
4:CE:36:THR:CG2	4:CE:37:VAL:H	2.25	0.45
5:CF:22:ILE:O	5:CF:26:THR:HG23	2.16	0.45
7:CH:17:GLN:HG3	7:CH:71:VAL:HG21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:CJ:40:ILE:CD1	9:CJ:73:LEU:HB3	2.46	0.45
11:CL:72:ASN:ND2	11:CL:73:LEU:H	2.15	0.45
13:CN:14:ALA:HA	13:CN:17:ASP:OD2	2.16	0.45
1:CA:564:C:C5	16:CQ:32:ILE:HD11	2.52	0.45
47:D0:43:THR:O	47:D0:44:ALA:HB3	2.16	0.45
22:DA:66:A:O2'	22:DA:67:G:C5'	2.61	0.45
23:DB:1199:U:C2	23:DB:1200:C:C6	3.05	0.45
23:DB:1790:C:H2'	23:DB:1791:A:C8	2.52	0.45
23:DB:2340:A:H2'	23:DB:2341:G:H8	1.81	0.45
23:DB:2363:G:O2'	23:DB:2364:C:H5'	2.16	0.45
23:DB:2684:U:H4'	32:DK:76:VAL:CG2	2.43	0.45
23:DB:2835:A:H61	23:DB:2878:U:H2'	1.80	0.45
23:DB:2886:A:C2	23:DB:2887:A:N7	2.85	0.45
23:DB:324:A:H61	23:DB:338:G:C2'	2.30	0.45
23:DB:544:C:O2'	23:DB:545:U:O4'	2.34	0.45
26:DD:23:PRO:HB3	26:DD:188:LEU:HG	1.97	0.45
23:DB:38:A:H1'	27:DE:43:THR:O	2.17	0.45
22:DA:57:A:H2	28:DF:29:ARG:HH22	1.61	0.45
28:DF:35:LEU:HD23	28:DF:153:ILE:HG12	1.99	0.45
30:DH:122:LEU:CD1	30:DH:146:VAL:HG22	2.41	0.45
23:DB:1098:A:C3'	52:DI:3:LYS:CA	2.69	0.45
52:DI:53:PRO:HG2	52:DI:77:VAL:HG11	1.98	0.45
33:DL:78:ARG:HA	33:DL:113:ALA:HB2	1.97	0.45
34:DM:81:ARG:HB2	34:DM:81:ARG:NH1	2.32	0.45
37:DP:26:GLU:CD	37:DP:46:VAL:HG13	2.37	0.45
39:DR:89:HIS:O	39:DR:90:ARG:CG	2.57	0.45
43:DW:47:GLY:HA2	43:DW:71:LYS:O	2.16	0.45
45:DY:6:ILE:HD11	45:DY:35:VAL:CB	2.47	0.45
1:AA:36:C:H2'	1:AA:37:U:O4'	2.17	0.45
1:AA:731:G:OP1	1:AA:766:A:H1'	2.17	0.45
1:AA:847:G:H2'	1:AA:848:C:H6	1.82	0.45
1:AA:89:U:O2'	1:AA:90:C:O5'	2.33	0.45
20:AB:186:VAL:CG2	20:AB:198:VAL:HG23	2.47	0.45
2:AC:156:LEU:CD1	2:AC:156:LEU:H	2.11	0.45
2:AC:185:THR:HG22	2:AC:186:SER:N	2.26	0.45
2:AC:8:GLY:HA2	2:AC:11:LEU:CG	2.43	0.45
3:AD:172:VAL:HA	3:AD:178:GLU:O	2.16	0.45
7:AH:54:THR:HG23	7:AH:55:LYS:HG2	1.99	0.45
9:AJ:87:LEU:H	9:AJ:88:MET:HE1	1.82	0.45
1:AA:796:C:H4'	10:AK:126:ARG:HH21	1.82	0.45
49:B2:18:PHE:HB3	49:B2:19:ARG:H	1.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:BA:94:A:H2'	22:BA:95:U:O4'	2.17	0.45
22:BA:99:A:H2'	22:BA:99:A:N3	2.32	0.45
23:BB:1275:A:N3	23:BB:1275:A:C3'	2.78	0.45
23:BB:340:A:H2'	23:BB:341:C:C5'	2.47	0.45
23:BB:350:G:H2'	23:BB:351:C:H6	1.82	0.45
23:BB:515:A:H3'	23:BB:516:C:H6	1.80	0.45
23:BB:829:A:H4'	55:BB:3324:HOH:O	2.16	0.45
25:BC:140:VAL:CG1	25:BC:163:ILE:HD11	2.46	0.45
26:BD:52:THR:HG22	26:BD:75:ALA:CB	2.44	0.45
27:BE:191:ASP:C	27:BE:193:VAL:H	2.20	0.45
23:BB:37:C:C2'	27:BE:46:GLN:CG	2.77	0.45
27:BE:57:LYS:HA	27:BE:57:LYS:CE	2.47	0.45
23:BB:2759:G:H21	29:BG:138:GLN:HE22	1.64	0.45
30:BH:24:GLY:HA2	30:BH:27:ARG:HH12	1.81	0.45
30:BH:39:ALA:O	30:BH:44:ILE:HD11	2.17	0.45
30:BH:94:ILE:HA	30:BH:98:ASP:OD2	2.17	0.45
31:BJ:93:ILE:CD1	31:BJ:100:VAL:HG11	2.46	0.45
32:BK:28:SER:O	32:BK:29:HIS:HB2	2.17	0.45
33:BL:122:VAL:HG23	33:BL:140:GLY:O	2.16	0.45
33:BL:60:ARG:HD2	50:B3:7:ARG:CZ	2.47	0.45
37:BP:114:ASN:N	37:BP:114:ASN:ND2	2.61	0.45
37:BP:15:ASP:N	37:BP:15:ASP:OD1	2.49	0.45
38:BQ:33:VAL:CG1	38:BQ:34:ALA:H	2.14	0.45
38:BQ:61:ILE:HB	38:BQ:75:TYR:CZ	2.52	0.45
39:BR:16:GLU:HG3	39:BR:17:GLY:H	1.82	0.45
40:BS:31:GLN:HB3	40:BS:35:ILE:HG22	1.99	0.45
41:BT:61:LEU:CD1	41:BT:62:VAL:H	2.28	0.45
42:BU:44:HIS:HB2	42:BU:56:GLY:H	1.80	0.45
45:BY:27:GLY:C	45:BY:28:LEU:HD12	2.37	0.45
45:BY:3:THR:O	45:BY:6:ILE:HG23	2.17	0.45
1:CA:1102:A:O2'	1:CA:1103:C:H5'	2.17	0.45
1:CA:1130:A:H2'	1:CA:1131:G:H8	1.81	0.45
1:CA:1453:G:H2'	1:CA:1454:G:O4'	2.17	0.45
1:CA:404:G:H2'	1:CA:405:U:C6	2.52	0.45
1:CA:551:U:O2'	1:CA:552:U:H5'	2.17	0.45
1:CA:611:C:H2'	1:CA:612:C:H6	1.80	0.45
1:CA:902:G:H2'	1:CA:903:G:C8	2.51	0.45
1:CA:1158:C:O2'	20:CB:131:LYS:HD3	2.16	0.45
2:CC:177:LEU:HD22	2:CC:177:LEU:H	1.81	0.45
2:CC:59:PRO:CG	2:CC:62:SER:HB2	2.33	0.45
3:CD:141:VAL:O	3:CD:141:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:62:MET:HG3	5:CF:64:VAL:CG2	2.42	0.45
8:CI:26:LYS:HG3	8:CI:61:ASP:HB2	1.97	0.45
10:CK:69:CYS:O	10:CK:73:VAL:N	2.49	0.45
1:CA:451:A:OP2	15:CP:70:ARG:NH2	2.50	0.45
18:CS:39:ILE:HB	18:CS:66:VAL:HA	1.98	0.45
48:D1:31:GLU:HG2	48:D1:32:LYS:HG3	1.98	0.45
23:DB:1175:A:H2'	23:DB:1176:U:C5'	2.42	0.45
23:DB:1700:A:C2'	23:DB:1701:A:H5'	2.44	0.45
23:DB:1760:C:H2'	23:DB:1761:C:O4'	2.16	0.45
23:DB:1773:A:C8	23:DB:1829:A:C8	3.05	0.45
23:DB:2259:U:C6	23:DB:2427:C:C5	3.05	0.45
23:DB:2420:C:C5	50:D3:30:HIS:O	2.70	0.45
23:DB:476:G:N2	23:DB:478:A:H3'	2.32	0.45
23:DB:483:A:C2'	23:DB:484:C:H5'	2.43	0.45
23:DB:633:A:H2'	23:DB:634:C:O4'	2.16	0.45
23:DB:747:U:O3'	40:DS:89:ALA:HB3	2.17	0.45
23:DB:844:A:C2	23:DB:845:A:N1	2.84	0.45
25:DC:163:ILE:HG22	25:DC:164:VAL:N	2.32	0.45
23:DB:730:A:C4'	25:DC:16:VAL:HG13	2.47	0.45
23:DB:1796:U:O3'	25:DC:251:THR:HA	2.16	0.45
25:DC:63:ILE:CG2	25:DC:64:VAL:N	2.79	0.45
27:DE:149:ILE:HD11	27:DE:187:VAL:N	2.31	0.45
29:DG:71:LEU:O	29:DG:74:MET:HB2	2.16	0.45
31:DJ:14:ASP:HB3	31:DJ:16:TYR:CD1	2.52	0.45
32:DK:71:ARG:HB3	32:DK:72:PRO:CD	2.46	0.45
33:DL:118:THR:HG23	33:DL:119:PRO:HD3	1.97	0.45
33:DL:39:LYS:NZ	33:DL:39:LYS:N	2.64	0.45
26:DD:186:LEU:HD11	37:DP:5:LYS:HE2	1.98	0.45
37:DP:83:ILE:H	37:DP:83:ILE:HG13	1.51	0.45
38:DQ:87:VAL:CG2	39:DR:54:VAL:HG11	2.46	0.45
39:DR:76:LYS:HA	39:DR:91:GLN:N	2.31	0.45
40:DS:10:ALA:O	40:DS:11:ARG:HB3	2.17	0.45
40:DS:1:MET:HB3	40:DS:4:ILE:HD12	1.98	0.45
42:DU:82:VAL:HB	42:DU:94:PHE:CB	2.43	0.45
36:DO:18:LEU:CD1	43:DW:76:ARG:HH21	2.29	0.45
44:DX:26:PHE:HA	44:DX:29:ARG:HD2	1.97	0.45
46:DZ:54:GLY:H	46:DZ:57:VAL:CG2	2.14	0.45
1:AA:1522:U:O2'	1:AA:1523:G:H5'	2.16	0.45
1:AA:238:A:C3'	1:AA:239:U:H5''	2.47	0.45
1:AA:404:G:H2'	1:AA:405:U:C6	2.52	0.45
1:AA:607:A:H2'	1:AA:608:A:C8	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:740:U:O3'	14:AO:38:LEU:HD21	2.16	0.45
1:AA:81:A:N3	1:AA:82:G:N7	2.65	0.45
20:AB:59:ILE:C	20:AB:59:ILE:HD12	2.37	0.45
2:AC:51:VAL:HA	2:AC:69:THR:HB	1.99	0.45
3:AD:99:ASN:HD21	3:AD:110:ARG:HE	1.64	0.45
3:AD:186:GLU:C	3:AD:190:LEU:HD13	2.37	0.45
6:AG:59:GLU:O	6:AG:63:VAL:HG23	2.17	0.45
9:AJ:68:ARG:HB3	9:AJ:70:HIS:NE2	2.32	0.45
9:AJ:87:LEU:HD13	9:AJ:87:LEU:C	2.37	0.45
11:AL:35:ARG:HG3	11:AL:36:VAL:N	2.27	0.45
12:AM:77:LYS:HE3	12:AM:81:ASP:HB3	1.99	0.45
15:AP:22:ALA:HB2	15:AP:32:PHE:HB3	1.98	0.45
47:B0:28:SER:HB2	47:B0:34:GLY:H	1.78	0.45
47:B0:42:ILE:HG21	47:B0:47:TYR:CE1	2.51	0.45
51:B4:13:ASN:ND2	51:B4:13:ASN:N	2.63	0.45
22:BA:109:A:H2'	22:BA:110:C:C6	2.51	0.45
22:BA:22:U:H2'	22:BA:23:G:C8	2.51	0.45
23:BB:1019:U:O2'	23:BB:1021:A:C2	2.64	0.45
23:BB:1057:A:C8	23:BB:1086:A:C8	3.05	0.45
23:BB:1184:U:H2'	23:BB:1185:G:C8	2.52	0.45
23:BB:1601:G:O2'	23:BB:1602:U:H5'	2.16	0.45
23:BB:2340:A:H2'	23:BB:2341:G:H8	1.82	0.45
23:BB:2772:C:H2'	23:BB:2773:C:H6	1.81	0.45
23:BB:2821:A:H2'	23:BB:2822:G:H8	1.82	0.45
23:BB:481:G:OP2	42:BU:43:LYS:HG3	2.16	0.45
23:BB:741:U:H2'	23:BB:742:A:H8	1.80	0.45
23:BB:85:G:H5''	42:BU:28:LEU:HD23	1.99	0.45
23:BB:997:G:O2'	23:BB:998:C:H5'	2.17	0.45
26:BD:13:ARG:HB2	26:BD:21:SER:CB	2.46	0.45
26:BD:26:VAL:O	26:BD:26:VAL:HG23	2.17	0.45
26:BD:70:LYS:HA	26:BD:70:LYS:CE	2.37	0.45
28:BF:148:VAL:O	28:BF:150:GLY:N	2.50	0.45
30:BH:121:VAL:HG21	30:BH:128:HIS:NE2	2.31	0.45
30:BH:21:VAL:HG13	30:BH:22:LYS:O	2.16	0.45
34:BM:68:PHE:HA	34:BM:69:PRO:HD2	1.76	0.45
35:BN:22:ARG:NH1	35:BN:69:ARG:HA	2.32	0.45
36:BO:41:ALA:H	36:BO:48:LEU:HB3	1.81	0.45
37:BP:105:LYS:HG3	37:BP:107:ALA:H	1.81	0.45
31:BJ:45:THR:HB	38:BQ:63:ARG:CZ	2.46	0.45
39:BR:62:GLU:OE1	39:BR:62:GLU:N	2.49	0.45
39:BR:35:PHE:HB2	39:BR:64:VAL:CG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BQ:39:ILE:CG2	39:BR:80:ARG:HD3	2.47	0.45
41:BT:29:THR:O	41:BT:30:ILE:HD12	2.17	0.45
41:BT:66:LYS:O	41:BT:67:VAL:HB	2.17	0.45
24:BV:72:VAL:HA	24:BV:94:ALA:H	1.81	0.45
24:BV:83:LYS:O	24:BV:85:LYS:N	2.49	0.45
41:BT:9:LYS:HG2	44:BX:22:LEU:HD11	1.99	0.45
45:BY:37:ARG:HH11	45:BY:37:ARG:HG3	1.82	0.45
1:CA:1136:C:H6	1:CA:1136:C:O5'	2.00	0.45
1:CA:1288:A:N1	1:CA:1371:G:H1'	2.31	0.45
1:CA:1387:G:H2'	1:CA:1388:C:H6	1.79	0.45
1:CA:1451:U:C5'	1:CA:1452:C:OP2	2.59	0.45
1:CA:1516:G:N1	1:CA:1519:A:OP2	2.50	0.45
1:CA:334:C:H2'	1:CA:335:C:H6	1.81	0.45
1:CA:975:A:OP2	1:CA:975:A:H4'	2.16	0.45
20:CB:34:ARG:NE	20:CB:39:ILE:HG13	2.32	0.45
2:CC:41:TYR:N	2:CC:41:TYR:CD2	2.83	0.45
4:CE:110:MET:O	4:CE:114:LEU:HG	2.16	0.45
8:CI:120:ALA:O	8:CI:121:ARG:HG2	2.17	0.45
11:CL:13:ARG:O	11:CL:14:LYS:HB3	2.16	0.45
11:CL:14:LYS:HZ3	11:CL:16:ALA:H	1.65	0.45
14:CO:69:LEU:HD12	14:CO:77:TYR:N	2.31	0.45
15:CP:48:GLU:HG2	15:CP:50:THR:H	1.81	0.45
21:CU:16:ARG:HA	21:CU:19:LYS:HZ3	1.82	0.45
21:CU:33:ARG:HB3	21:CU:34:ARG:H	1.65	0.45
48:D1:8:ILE:CG2	48:D1:27:ARG:HD3	2.46	0.45
51:D4:10:LEU:HD12	51:D4:10:LEU:N	2.32	0.45
23:DB:1076:C:O2'	23:DB:1077:A:H5'	2.17	0.45
23:DB:1791:A:N1	23:DB:1829:A:H5'	2.31	0.45
23:DB:1803:A:H2	23:DB:1823:G:C1'	2.29	0.45
23:DB:1821:A:H5'	25:DC:155:ARG:NH2	2.28	0.45
23:DB:1843:C:H2'	23:DB:1844:C:H6	1.81	0.45
23:DB:2188:U:H2'	23:DB:2189:U:O4'	2.16	0.45
23:DB:2464:G:H2'	23:DB:2465:C:C6	2.52	0.45
23:DB:257:C:H3'	23:DB:258:G:H8	1.82	0.45
23:DB:259:G:O2'	23:DB:260:G:H5'	2.16	0.45
23:DB:2685:G:OP1	37:DP:62:LYS:NZ	2.47	0.45
23:DB:2884:U:C2	47:D0:48:TYR:CZ	3.04	0.45
23:DB:38:A:N3	27:DE:43:THR:HG22	2.32	0.45
23:DB:40:U:H2'	23:DB:41:C:C6	2.52	0.45
23:DB:437:U:O2'	23:DB:438:G:H5'	2.16	0.45
23:DB:904:G:H2'	23:DB:905:A:H8	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:99:U:O2	23:DB:99:U:O4'	2.33	0.45
25:DC:130:PRO:HA	25:DC:188:ARG:HA	1.98	0.45
25:DC:156:SER:O	25:DC:194:VAL:O	2.35	0.45
27:DE:135:ALA:O	27:DE:139:LYS:HB3	2.16	0.45
52:DI:89:SER:HA	52:DI:97:VAL:CG1	2.47	0.45
38:DQ:101:ASP:OD1	38:DQ:104:ALA:HB3	2.17	0.45
39:DR:44:GLY:N	39:DR:53:PHE:CE2	2.84	0.45
39:DR:69:GLY:HA2	39:DR:97:LYS:N	2.29	0.45
42:DU:3:LYS:O	42:DU:4:ILE:HB	2.17	0.45
1:AA:1399:C:H1'	55:AA:1970:HOH:O	2.16	0.45
1:AA:1515:G:O2'	1:AA:1516:G:H5'	2.17	0.45
1:AA:24:U:O2'	1:AA:25:C:H5'	2.16	0.45
1:AA:402:G:H2'	1:AA:403:C:C6	2.52	0.45
1:AA:662:U:H2'	1:AA:663:A:H8	1.76	0.45
20:AB:140:LEU:HB3	20:AB:144:GLU:OE2	2.17	0.45
2:AC:153:SER:CB	2:AC:196:GLY:H	2.27	0.45
2:AC:5:HIS:O	2:AC:9:ILE:HG22	2.16	0.45
8:AI:51:LEU:HD23	8:AI:56:MET:HE3	1.99	0.45
11:AL:17:LYS:HB2	11:AL:17:LYS:HE3	1.82	0.45
12:AM:8:ILE:O	12:AM:8:ILE:HG13	2.17	0.45
13:AN:81:ILE:HD12	13:AN:82:LYS:N	2.31	0.45
16:AQ:74:LEU:HD13	16:AQ:74:LEU:C	2.37	0.45
18:AS:29:PRO:CA	18:AS:47:THR:HB	2.47	0.45
50:B3:35:LYS:O	50:B3:36:ALA:HB3	2.16	0.45
22:BA:78:A:O2'	22:BA:79:G:H5'	2.17	0.45
23:BB:1100:C:H2'	23:BB:1101:U:H6	1.81	0.45
23:BB:1322:A:H2'	23:BB:1323:C:H5'	1.98	0.45
23:BB:1509:A:C5'	23:BB:1510:G:H5'	2.47	0.45
23:BB:1683:U:H2'	23:BB:1684:G:C8	2.51	0.45
23:BB:1987:A:H2'	23:BB:1988:G:H8	1.82	0.45
23:BB:2465:C:O2'	23:BB:2466:C:H5'	2.16	0.45
23:BB:2708:G:H4'	35:BN:68:ALA:HB1	1.99	0.45
23:BB:2830:C:O2'	23:BB:2831:G:H5'	2.17	0.45
23:BB:992:C:H2'	23:BB:993:G:C8	2.51	0.45
25:BC:153:LEU:CD2	25:BC:153:LEU:H	2.26	0.45
25:BC:224:MET:HE2	25:BC:233:GLY:O	2.17	0.45
26:BD:128:ARG:NH1	26:BD:144:GLY:HA2	2.32	0.45
27:BE:7:ASP:O	27:BE:8:ALA:HB3	2.17	0.45
30:BH:2:GLN:HE22	30:BH:20:ASN:N	2.14	0.45
52:BI:38:CYS:O	52:BI:42:ASN:ND2	2.49	0.45
35:BN:73:ASN:HA	35:BN:76:VAL:CG1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BQ:103:VAL:HG23	38:BQ:104:ALA:N	2.31	0.45
38:BQ:73:ILE:HG13	38:BQ:77:LYS:HD2	1.99	0.45
23:BB:139:U:C2	41:BT:1:MET:HG2	2.51	0.45
43:BW:24:ARG:HB3	43:BW:57:THR:O	2.16	0.45
1:CA:1098:C:O2'	1:CA:1099:G:H5'	2.16	0.45
1:CA:1128:C:H4'	1:CA:1148:U:O2	2.17	0.45
1:CA:1382:C:H5'	6:CG:78:ARG:HE	1.82	0.45
1:CA:1488:G:H2'	1:CA:1489:G:H8	1.82	0.45
1:CA:239:U:H6	1:CA:239:U:H5'	1.80	0.45
20:CB:211:LEU:HG	20:CB:212:TYR:N	2.32	0.45
2:CC:10:ARG:HG2	2:CC:10:ARG:HH11	1.82	0.45
4:CE:82:HIS:NE2	4:CE:146:MET:HB2	2.31	0.45
4:CE:89:THR:C	4:CE:91:SER:H	2.18	0.45
7:CH:29:SER:O	7:CH:30:LYS:C	2.55	0.45
11:CL:82:ARG:HB2	11:CL:97:VAL:HG13	1.98	0.45
1:CA:1227:A:H4'	12:CM:113:LYS:HE2	1.97	0.45
14:CO:63:ARG:HH21	14:CO:87:ARG:NH1	2.15	0.45
15:CP:32:PHE:HD1	15:CP:32:PHE:C	2.20	0.45
16:CQ:60:ILE:HG22	16:CQ:74:LEU:HA	1.98	0.45
22:DA:6:G:O2'	22:DA:7:G:H5'	2.16	0.45
23:DB:1286:A:H1'	23:DB:1288:G:OP2	2.17	0.45
23:DB:2144:G:C2	23:DB:2146:C:H5'	2.52	0.45
23:DB:2651:C:O2'	23:DB:2652:C:H5'	2.17	0.45
23:DB:2673:G:H2'	23:DB:2674:G:C8	2.52	0.45
23:DB:510:C:O2'	23:DB:1236:G:H5'	2.17	0.45
23:DB:544:C:H2'	23:DB:545:U:C2	2.52	0.45
23:DB:547:A:N3	23:DB:547:A:C2'	2.77	0.45
23:DB:97:C:H2'	23:DB:98:G:O4'	2.17	0.45
23:DB:1820:U:N3	25:DC:197:ALA:HB1	2.27	0.45
25:DC:259:ASN:O	25:DC:261:ARG:HG3	2.17	0.45
25:DC:86:ARG:HH11	25:DC:86:ARG:HB2	1.81	0.45
28:DF:7:TYR:HA	28:DF:11:VAL:HB	1.98	0.45
29:DG:103:ASN:HA	29:DG:112:VAL:O	2.17	0.45
29:DG:100:ASN:HA	29:DG:116:LEU:HD11	1.99	0.45
29:DG:141:GLY:O	29:DG:144:ALA:HB3	2.16	0.45
52:DI:99:LYS:HB2	52:DI:140:GLU:OE1	2.16	0.45
31:DJ:73:VAL:HA	31:DJ:88:THR:O	2.17	0.45
33:DL:92:LEU:HB2	33:DL:125:LEU:HD11	1.99	0.45
34:DM:33:LEU:HD21	34:DM:124:LEU:HD13	1.98	0.45
36:DO:98:GLN:O	36:DO:99:TYR:HB2	2.17	0.45
40:DS:36:LEU:HA	40:DS:39:THR:CG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:49:LYS:O	41:DT:51:PHE:N	2.41	0.45
42:DU:27:VAL:CB	42:DU:33:VAL:HG22	2.46	0.45
44:DX:17:GLU:HA	44:DX:21:LEU:CB	2.46	0.45
1:AA:1096:C:H2'	1:AA:1097:C:C6	2.52	0.45
1:AA:1250:A:O2'	1:AA:1251:A:H5'	2.17	0.45
1:AA:295:C:H2'	1:AA:296:U:C6	2.52	0.45
1:AA:408:A:OP1	3:AD:111:ALA:HB3	2.17	0.45
1:AA:202:G:N2	1:AA:465:A:H61	2.14	0.45
20:AB:96:LEU:H	20:AB:99:MET:HE3	1.82	0.45
2:AC:135:ARG:C	2:AC:137:VAL:H	2.20	0.45
1:AA:543:U:P	3:AD:13:ARG:HH21	2.40	0.45
5:AF:66:ALA:HB1	5:AF:67:PRO:HD2	1.99	0.45
1:AA:1380:U:C4	6:AG:2:ARG:HG3	2.52	0.45
10:AK:80:ASN:HD21	10:AK:107:THR:HG23	1.82	0.45
11:AL:30:ARG:HH11	11:AL:30:ARG:CB	2.29	0.45
11:AL:52:CYS:SG	11:AL:66:ILE:HD11	2.56	0.45
12:AM:44:ILE:CA	12:AM:47:LEU:HD13	2.47	0.45
13:AN:60:ARG:HE	13:AN:62:ARG:CZ	2.30	0.45
15:AP:4:ILE:HA	15:AP:20:VAL:O	2.16	0.45
16:AQ:23:ALA:HB1	16:AQ:40:THR:CG2	2.47	0.45
33:BL:60:ARG:HD3	50:B3:11:LYS:NZ	2.31	0.45
50:B3:44:ARG:HG2	50:B3:44:ARG:NH1	2.32	0.45
23:BB:1309:G:H4'	49:B2:7:PRO:CG	2.45	0.45
23:BB:143:C:C3'	23:BB:144:A:C8	3.00	0.45
23:BB:1547:C:H2'	23:BB:1548:A:C8	2.52	0.45
23:BB:15:G:H4'	47:B0:17:SER:OG	2.17	0.45
23:BB:1669:A:H2'	23:BB:1669:A:N3	2.30	0.45
23:BB:2291:U:H2'	23:BB:2292:U:H6	1.80	0.45
23:BB:2319:G:C2	23:BB:2321:U:O4	2.70	0.45
23:BB:2651:C:O2'	23:BB:2652:C:H5'	2.17	0.45
23:BB:2673:G:H2'	23:BB:2674:G:C8	2.52	0.45
23:BB:2776:A:H4'	23:BB:2777:G:C5'	2.46	0.45
23:BB:1752:C:H4'	23:BB:2862:G:OP1	2.16	0.45
23:BB:338:G:C2	23:BB:339:U:C1'	2.99	0.45
23:BB:402:A:H2'	23:BB:403:U:O4'	2.16	0.45
23:BB:463:G:N2	23:BB:466:A:OP2	2.48	0.45
23:BB:527:C:H5'	55:BB:3419:HOH:O	2.17	0.45
25:BC:208:GLY:C	25:BC:210:ALA:H	2.20	0.45
26:BD:148:GLN:HB3	26:BD:151:THR:HG21	1.98	0.45
26:BD:38:LYS:HG2	26:BD:39:ASP:N	2.22	0.45
23:BB:1255:U:H2'	27:BE:67:ARG:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:135:ILE:HG23	28:BF:136:ILE:N	2.31	0.45
28:BF:31:GLU:HB3	28:BF:32:LYS:H	1.54	0.45
29:BG:120:ILE:HD12	29:BG:120:ILE:N	2.32	0.45
29:BG:121:THR:O	29:BG:133:LYS:HD3	2.16	0.45
30:BH:11:ASN:HD22	30:BH:12:LEU:N	2.10	0.45
52:BI:14:ALA:CB	52:BI:50:LYS:HA	2.47	0.45
32:BK:10:VAL:HG13	32:BK:19:VAL:CG2	2.46	0.45
32:BK:8:LEU:HB3	32:BK:84:CYS:SG	2.57	0.45
23:BB:811:U:C6	33:BL:29:LYS:HD3	2.52	0.45
33:BL:33:ARG:NH1	33:BL:34:GLY:HA3	2.32	0.45
34:BM:109:PRO:O	34:BM:112:LEU:HB2	2.17	0.45
35:BN:101:GLY:H	35:BN:110:MET:H	1.65	0.45
37:BP:38:ARG:CZ	37:BP:39:LEU:H	2.29	0.45
38:BQ:16:ILE:HD11	38:BQ:31:TYR:CD1	2.52	0.45
38:BQ:35:PHE:CA	38:BQ:39:ILE:HD12	2.46	0.45
1:CA:1319:A:H2'	1:CA:1323:G:N7	2.32	0.45
1:CA:224:U:H2'	1:CA:225:C:H6	1.82	0.45
1:CA:295:C:H2'	1:CA:296:U:C6	2.52	0.45
1:CA:343:U:H2'	1:CA:345:C:C5	2.52	0.45
1:CA:432:A:C2'	1:CA:433:G:H5'	2.47	0.45
1:CA:478:A:H2'	1:CA:479:U:O4'	2.15	0.45
1:CA:957:U:H2'	1:CA:959:A:OP2	2.17	0.45
2:CC:53:ARG:HA	2:CC:113:LYS:CE	2.47	0.45
2:CC:17:TRP:C	2:CC:19:SER:H	2.20	0.45
2:CC:63:ILE:HG12	2:CC:65:VAL:CG2	2.46	0.45
3:CD:104:MET:HG3	3:CD:142:VAL:HG21	1.99	0.45
3:CD:98:ASP:OD2	3:CD:99:ASN:N	2.50	0.45
4:CE:137:ARG:NH1	4:CE:137:ARG:HG2	2.31	0.45
4:CE:14:LEU:HD22	4:CE:15:ILE:N	2.32	0.45
4:CE:68:ARG:O	4:CE:69:ASN:HB3	2.17	0.45
8:CI:115:VAL:HG22	8:CI:116:GLY:N	2.32	0.45
1:CA:1250:A:C5'	8:CI:69:GLY:H	2.29	0.45
8:CI:88:GLU:HA	8:CI:91:GLU:OE2	2.16	0.45
11:CL:41:PRO:HG2	11:CL:45:ASN:O	2.16	0.45
17:CR:33:THR:HG23	17:CR:35:SER:H	1.81	0.45
18:CS:50:VAL:CG2	18:CS:70:LEU:HG	2.45	0.45
21:CU:7:GLU:OE1	21:CU:11:PHE:HZ	2.00	0.45
23:DB:1148:U:O2'	23:DB:1149:G:H5'	2.15	0.45
23:DB:1221:C:H2'	23:DB:1222:U:H6	1.82	0.45
23:DB:154:U:H2'	23:DB:155:A:C8	2.52	0.45
23:DB:1880:U:H2'	23:DB:1881:C:C6	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2065:C:H2'	23:DB:2066:C:C6	2.51	0.45
23:DB:2460:U:H2'	23:DB:2461:A:H8	1.81	0.45
23:DB:2649:C:H2'	23:DB:2650:U:C6	2.52	0.45
23:DB:2712:C:H2'	23:DB:2714:G:O3'	2.17	0.45
23:DB:2769:U:O2'	23:DB:2770:G:H5'	2.16	0.45
23:DB:303:G:H2'	23:DB:304:U:H6	1.82	0.45
23:DB:438:G:O2'	23:DB:439:A:H5'	2.16	0.45
25:DC:71:ASP:HA	25:DC:117:SER:OG	2.17	0.45
26:DD:14:ILE:O	26:DD:15:PHE:HD1	2.00	0.45
26:DD:172:VAL:CG1	26:DD:175:LEU:HD11	2.46	0.45
26:DD:77:ARG:HB2	26:DD:77:ARG:CZ	2.47	0.45
26:DD:7:LYS:HB3	26:DD:201:LEU:HD22	1.98	0.45
27:DE:17:THR:CG2	27:DE:18:THR:H	2.28	0.45
28:DF:15:LEU:HD13	28:DF:28:PRO:HG2	1.98	0.45
22:DA:55:U:O2'	28:DF:25:MET:HB3	2.16	0.45
29:DG:138:GLN:HE22	29:DG:142:GLN:HG3	1.82	0.45
30:DH:94:ILE:O	30:DH:122:LEU:HB2	2.16	0.45
30:DH:4:ILE:HD13	30:DH:37:VAL:HG13	1.98	0.45
30:DH:8:LYS:HA	30:DH:13:GLY:O	2.17	0.45
31:DJ:5:THR:HG23	31:DJ:7:LYS:HD3	1.99	0.45
33:DL:7:SER:CB	33:DL:8:PRO:CD	2.87	0.45
33:DL:94:THR:OG1	33:DL:95:LEU:HD12	2.16	0.45
34:DM:134:THR:HG22	34:DM:135:VAL:N	2.24	0.45
23:DB:1454:C:H5'	35:DN:63:ARG:HD2	1.97	0.45
36:DO:45:SER:C	36:DO:47:VAL:H	2.20	0.45
36:DO:86:GLY:O	36:DO:88:LYS:N	2.50	0.45
32:DK:78:ARG:H	37:DP:72:VAL:HG23	1.82	0.45
38:DQ:17:LEU:O	38:DQ:20:ALA:HB3	2.17	0.45
23:DB:1251:C:H5''	38:DQ:5:ARG:NH1	2.32	0.45
39:DR:11:GLN:N	39:DR:21:ARG:NH2	2.64	0.45
42:DU:85:ARG:O	42:DU:87:GLU:N	2.50	0.45
45:DY:3:THR:O	45:DY:36:GLU:HA	2.16	0.45
1:AA:1323:G:O2'	1:AA:1362:A:O4'	2.35	0.45
1:AA:1478:U:H2'	1:AA:1479:C:C6	2.52	0.45
1:AA:1481:U:H2'	1:AA:1482:G:C8	2.52	0.45
1:AA:1521:C:O2'	1:AA:1522:U:H5'	2.17	0.45
1:AA:385:C:O2'	1:AA:386:C:H5'	2.17	0.45
1:AA:408:A:P	3:AD:109:THR:HG21	2.57	0.45
1:AA:521:G:O2'	1:AA:522:C:H5'	2.17	0.45
1:AA:551:U:O2'	1:AA:552:U:H5'	2.17	0.45
1:AA:781:A:H2'	1:AA:782:A:C5'	2.38	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:957:U:H2'	1:AA:959:A:OP2	2.17	0.45
20:AB:23:ASN:HD22	20:AB:23:ASN:C	2.20	0.45
20:AB:81:ASP:O	20:AB:85:SER:HB2	2.17	0.45
3:AD:145:ARG:HH11	3:AD:145:ARG:CB	2.29	0.45
3:AD:94:GLU:HA	3:AD:94:GLU:OE2	2.16	0.45
4:AE:56:PRO:O	4:AE:59:ILE:HG22	2.16	0.45
6:AG:73:GLU:OE2	6:AG:90:VAL:HG12	2.17	0.45
1:AA:1148:U:H5'	8:AI:6:TYR:OH	2.17	0.45
12:AM:52:ILE:HG23	12:AM:53:ASP:N	2.32	0.45
14:AO:77:TYR:O	14:AO:81:ILE:HG23	2.17	0.45
16:AQ:29:LYS:HD3	16:AQ:35:LYS:CA	2.47	0.45
17:AR:22:TYR:CD2	17:AR:60:ARG:HD2	2.51	0.45
18:AS:64:GLU:O	18:AS:66:VAL:HG23	2.17	0.45
48:B1:20:TYR:O	48:B1:21:THR:HB	2.17	0.45
48:B1:40:PRO:HA	48:B1:48:TYR:OH	2.16	0.45
50:B3:29:ARG:HA	50:B3:33:THR:HG21	1.95	0.45
23:BB:640:C:H1'	50:B3:44:ARG:NH2	2.32	0.45
51:B4:24:ARG:HH22	51:B4:36:ARG:HD3	1.79	0.45
51:B4:7:VAL:HB	51:B4:35:GLN:NE2	2.31	0.45
23:BB:1064:C:C4'	52:BI:90:GLY:HA2	2.47	0.45
23:BB:1487:U:H2'	23:BB:1488:C:C6	2.50	0.45
23:BB:1736:U:H2'	23:BB:1737:G:C8	2.52	0.45
23:BB:1985:C:O2'	23:BB:1986:C:H5'	2.17	0.45
23:BB:2065:C:H2'	23:BB:2066:C:C6	2.51	0.45
23:BB:2671:G:H2'	23:BB:2672:U:H6	1.80	0.45
23:BB:493:G:H2'	23:BB:494:G:O4'	2.17	0.45
23:BB:952:G:C6	23:BB:966:G:C6	3.05	0.45
26:BD:128:ARG:CA	26:BD:128:ARG:NE	2.80	0.45
28:BF:132:ARG:NH2	28:BF:147:ARG:HE	2.15	0.45
28:BF:13:LYS:HD3	28:BF:16:MET:HG3	1.99	0.45
32:BK:20:MET:O	32:BK:42:THR:N	2.49	0.45
33:BL:77:ILE:HD12	33:BL:110:VAL:HG22	1.99	0.45
34:BM:31:PHE:HA	34:BM:128:THR:HG22	1.97	0.45
34:BM:7:THR:C	34:BM:8:LYS:HD3	2.37	0.45
35:BN:38:LEU:N	35:BN:39:PRO:CD	2.80	0.45
37:BP:26:GLU:O	37:BP:46:VAL:HG22	2.17	0.45
37:BP:9:GLN:HB3	37:BP:56:SER:HB2	1.97	0.45
38:BQ:16:ILE:HD11	38:BQ:31:TYR:HD1	1.82	0.45
38:BQ:19:GLN:NE2	38:BQ:38:VAL:HG23	2.32	0.45
38:BQ:92:LYS:CD	38:BQ:93:ILE:N	2.80	0.45
40:BS:70:LYS:O	40:BS:71:VAL:HB	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:2:ALA:O	42:BU:3:LYS:HD3	2.16	0.45
1:CA:1147:C:HO2'	8:CI:6:TYR:HE2	1.64	0.45
1:CA:24:U:O2'	1:CA:25:C:H5'	2.16	0.45
1:CA:285:C:H2'	1:CA:286:C:H6	1.81	0.45
1:CA:817:C:C2	1:CA:819:A:O4'	2.70	0.45
1:CA:953:G:H3'	1:CA:954:G:H8	1.82	0.45
2:CC:129:PHE:CG	2:CC:130:ARG:N	2.85	0.45
6:CG:29:LEU:HG	6:CG:104:VAL:HG22	1.99	0.45
11:CL:89:LEU:N	11:CL:89:LEU:HD22	2.32	0.45
11:CL:91:GLY:O	11:CL:93:ARG:NE	2.50	0.45
12:CM:73:SER:O	12:CM:77:LYS:HG3	2.17	0.45
15:CP:57:ILE:HD11	15:CP:75:ILE:HD11	1.98	0.45
47:D0:29:VAL:HG23	47:D0:32:THR:OG1	2.16	0.45
48:D1:14:ALA:CB	48:D1:19:PHE:HB3	2.47	0.45
23:DB:2420:C:OP2	50:D3:31:ILE:HA	2.17	0.45
22:DA:13:G:H2'	22:DA:14:U:H5''	1.99	0.45
23:DB:1190:G:OP1	33:DL:38:GLN:CB	2.65	0.45
23:DB:1509:A:C5'	23:DB:1510:G:H5'	2.47	0.45
23:DB:1585:C:H2'	23:DB:1586:A:C8	2.52	0.45
23:DB:1712:U:H3'	23:DB:1713:A:C5'	2.46	0.45
23:DB:1131:G:N7	23:DB:2025:C:H4'	2.32	0.45
23:DB:2637:U:O2'	23:DB:2638:G:H5'	2.17	0.45
23:DB:2733:A:H2'	23:DB:2734:A:O4'	2.17	0.45
23:DB:743:A:OP1	26:DD:134:HIS:CD2	2.70	0.45
23:DB:743:A:OP1	26:DD:134:HIS:NE2	2.50	0.45
26:DD:40:LEU:O	26:DD:41:ALA:HB3	2.17	0.45
23:DB:1205:A:N7	27:DE:164:LEU:CD1	2.80	0.45
30:DH:4:ILE:O	30:DH:5:LEU:HD22	2.17	0.45
30:DH:7:ASP:HB3	30:DH:8:LYS:H	1.60	0.45
31:DJ:58:ASN:O	31:DJ:60:ASP:N	2.45	0.45
31:DJ:94:ALA:HB3	31:DJ:95:ARG:HE	1.82	0.45
35:DN:45:ARG:NH2	35:DN:97:ILE:HG12	2.31	0.45
39:DR:37:GLU:CG	39:DR:62:GLU:HG3	2.45	0.45
41:DT:15:HIS:HB3	41:DT:31:VAL:CG1	2.47	0.45
41:DT:85:VAL:HG23	41:DT:86:THR:N	2.32	0.45
1:AA:1282:C:H2'	1:AA:1283:U:C6	2.52	0.44
1:AA:1439:G:H2'	1:AA:1440:U:O4'	2.18	0.44
1:AA:432:A:C2'	1:AA:433:G:H5'	2.47	0.44
1:AA:649:A:H2'	1:AA:650:G:O4'	2.17	0.44
1:AA:817:C:C2	1:AA:819:A:O4'	2.70	0.44
2:AC:25:THR:CG2	13:AN:74:ARG:HD2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:10:LEU:HD21	3:AD:62:ARG:CG	2.46	0.44
5:AF:40:GLU:CG	5:AF:42:TRP:HE1	2.27	0.44
13:AN:50:LEU:CG	13:AN:51:PRO:HD3	2.46	0.44
14:AO:70:LYS:HG2	14:AO:77:TYR:CG	2.52	0.44
15:AP:39:PHE:HE2	15:AP:70:ARG:NH2	2.15	0.44
17:AR:64:LEU:CB	17:AR:66:LEU:HG	2.47	0.44
18:AS:50:VAL:H	18:AS:57:VAL:H	1.64	0.44
21:AU:17:ARG:CA	21:AU:20:ARG:HB3	2.43	0.44
47:B0:51:ARG:O	47:B0:52:LYS:HD2	2.18	0.44
23:BB:2285:C:OP1	48:B1:26:LYS:HE3	2.17	0.44
23:BB:126:A:OP2	49:B2:45:SER:OG	2.29	0.44
23:BB:1292:G:H2'	23:BB:1293:C:H6	1.82	0.44
23:BB:1341:G:H2'	23:BB:1397:U:HO2'	1.81	0.44
23:BB:138:U:H2'	23:BB:140:C:C1'	2.45	0.44
23:BB:1773:A:C8	23:BB:1829:A:C8	3.05	0.44
23:BB:1880:U:H2'	23:BB:1881:C:C6	2.52	0.44
23:BB:2304:G:N2	23:BB:2312:U:H3	2.15	0.44
23:BB:2373:G:H2'	23:BB:2374:C:H6	1.80	0.44
23:BB:2464:G:H2'	23:BB:2465:C:C6	2.51	0.44
23:BB:2576:G:P	26:BD:149:ASN:CB	2.94	0.44
23:BB:699:A:H2'	23:BB:700:G:O4'	2.17	0.44
23:BB:807:U:H2'	23:BB:808:G:H8	1.81	0.44
23:BB:844:A:H2	23:BB:845:A:N1	2.16	0.44
23:BB:1790:C:H4'	25:BC:207:ALA:CB	2.46	0.44
25:BC:90:ILE:CG2	25:BC:91:ALA:N	2.79	0.44
23:BB:2774:C:P	26:BD:169:ARG:HG3	2.57	0.44
28:BF:101:ARG:HD2	28:BF:105:ILE:HD12	1.99	0.44
29:BG:158:GLY:CA	29:BG:170:THR:HG21	2.46	0.44
30:BH:134:VAL:HG22	30:BH:135:HIS:N	2.32	0.44
30:BH:147:VAL:HG12	30:BH:148:ALA:N	2.32	0.44
30:BH:35:LYS:HD2	30:BH:35:LYS:H	1.82	0.44
32:BK:24:VAL:HG12	32:BK:30:ARG:HB3	1.98	0.44
35:BN:86:ARG:HH21	35:BN:116:VAL:HG11	1.82	0.44
37:BP:31:VAL:HA	37:BP:81:ASP:CB	2.48	0.44
39:BR:6:GLN:O	39:BR:7:SER:OG	2.33	0.44
46:BZ:31:ASP:HA	46:BZ:47:LYS:O	2.17	0.44
1:CA:1426:G:H2'	1:CA:1427:C:C6	2.52	0.44
1:CA:1526:G:OP1	21:CU:37:TYR:HB3	2.17	0.44
1:CA:385:C:O2'	1:CA:386:C:H5'	2.17	0.44
3:CD:123:MET:CB	3:CD:128:VAL:HA	2.46	0.44
3:CD:123:MET:HB3	3:CD:128:VAL:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:172:VAL:HG22	3:CD:173:ASP:N	2.32	0.44
3:CD:18:LEU:HD22	3:CD:18:LEU:N	2.31	0.44
3:CD:74:TYR:HB3	3:CD:89:LEU:HD12	1.98	0.44
3:CD:71:PHE:CE2	3:CD:89:LEU:HD11	2.51	0.44
5:CF:52:ASN:O	5:CF:53:LYS:HB3	2.17	0.44
1:CA:1371:G:O3'	8:CI:70:GLY:HA3	2.17	0.44
9:CJ:19:ASP:HA	9:CJ:22:THR:HG22	1.99	0.44
10:CK:70:ALA:CA	10:CK:73:VAL:HG22	2.47	0.44
12:CM:84:CYS:O	12:CM:88:LEU:HG	2.16	0.44
50:D3:21:PHE:O	50:D3:48:MET:HB2	2.16	0.44
51:D4:9:LYS:O	51:D4:25:VAL:HA	2.18	0.44
22:DA:107:G:O2'	22:DA:108:A:H5'	2.17	0.44
23:DB:1005:C:O2'	23:DB:1006:C:H5'	2.15	0.44
23:DB:1025:G:H8	23:DB:1025:G:OP1	2.00	0.44
23:DB:1116:G:O2'	23:DB:1117:C:H5'	2.17	0.44
23:DB:1021:A:N6	23:DB:1142:A:N6	2.54	0.44
23:DB:1180:U:H2'	23:DB:1181:U:O4'	2.17	0.44
23:DB:1914:C:H3'	23:DB:1914:C:OP2	2.18	0.44
23:DB:231:A:H3'	23:DB:232:G:C8	2.52	0.44
23:DB:2457:U:O2'	23:DB:2458:G:H5'	2.17	0.44
23:DB:2598:A:P	25:DC:237:ARG:HH12	2.39	0.44
26:DD:45:TYR:HB3	26:DD:46:ARG:H	1.65	0.44
26:DD:31:ALA:HB3	26:DD:95:SER:HB2	1.99	0.44
28:DF:121:PHE:CE1	28:DF:166:ARG:HG2	2.52	0.44
30:DH:19:VAL:HG22	30:DH:20:ASN:N	2.32	0.44
30:DH:53:GLU:CD	30:DH:57:LYS:HD2	2.37	0.44
31:DJ:73:VAL:HG22	31:DJ:74:TYR:H	1.80	0.44
23:DB:1199:U:O2'	38:DQ:2:ARG:HB3	2.16	0.44
38:DQ:52:ARG:HG2	38:DQ:53:LYS:H	1.83	0.44
43:DW:76:ARG:C	43:DW:78:PHE:H	2.20	0.44
45:DY:17:PRO:HD2	45:DY:18:LYS:NZ	2.31	0.44
1:AA:1262:C:H2'	1:AA:1263:C:O4'	2.17	0.44
1:AA:178:C:O2'	1:AA:179:A:H5'	2.17	0.44
1:AA:343:U:H2'	1:AA:345:C:C5	2.52	0.44
1:AA:668:G:O2'	1:AA:669:G:H5'	2.17	0.44
1:AA:734:G:H2'	1:AA:735:C:C6	2.53	0.44
20:AB:67:LEU:HD12	20:AB:153:MET:CE	2.38	0.44
4:AE:108:GLY:H	4:AE:110:MET:CE	2.30	0.44
4:AE:16:ALA:O	4:AE:34:ALA:HA	2.16	0.44
6:AG:87:PRO:CG	6:AG:151:ALA:HB2	2.46	0.44
6:AG:74:VAL:HG12	6:AG:75:LYS:N	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:32:VAL:O	11:AL:33:CYS:HB2	2.15	0.44
13:AN:60:ARG:NE	13:AN:62:ARG:HE	2.15	0.44
15:AP:40:ASN:HB3	15:AP:49:GLY:O	2.17	0.44
15:AP:39:PHE:CE1	15:AP:74:LEU:HD22	2.52	0.44
16:AQ:44:HIS:O	16:AQ:72:TRP:HB2	2.16	0.44
23:BB:1263:U:H4'	47:B0:8:THR:HB	2.00	0.44
23:BB:1286:A:H1'	23:BB:1288:G:OP2	2.17	0.44
23:BB:1541:C:O2'	23:BB:1542:U:H5'	2.17	0.44
23:BB:1589:U:H2'	23:BB:1590:A:C8	2.52	0.44
23:BB:173:A:H2'	23:BB:174:U:C6	2.52	0.44
23:BB:2065:C:C2	23:BB:2066:C:C5	3.05	0.44
23:BB:2204:G:O2'	23:BB:2205:A:H5'	2.17	0.44
23:BB:2220:U:O2'	23:BB:2221:G:H5'	2.17	0.44
23:BB:2257:U:OP1	43:BW:5:ALA:HB2	2.18	0.44
23:BB:2354:C:H4'	43:BW:30:VAL:HG13	1.98	0.44
23:BB:2466:C:O2'	23:BB:2467:C:H5'	2.17	0.44
23:BB:2521:C:O2'	23:BB:2522:U:H5'	2.17	0.44
23:BB:2678:C:O2'	23:BB:2679:A:H5'	2.17	0.44
23:BB:269:C:O2'	23:BB:270:A:H5'	2.17	0.44
23:BB:273:G:H2'	23:BB:274:C:C6	2.53	0.44
23:BB:689:A:H2'	23:BB:690:G:C8	2.52	0.44
23:BB:701:G:O2'	23:BB:702:U:H5'	2.16	0.44
23:BB:718:A:H2'	23:BB:719:C:H5'	1.99	0.44
25:BC:115:ILE:HA	25:BC:126:GLY:CA	2.47	0.44
25:BC:27:LYS:HA	25:BC:79:ARG:CZ	2.48	0.44
26:BD:151:THR:HB	26:BD:152:PRO:CD	2.48	0.44
23:BB:2306:C:N4	28:BF:38:GLY:H	2.03	0.44
29:BG:154:GLU:OE2	29:BG:159:LYS:HG2	2.17	0.44
30:BH:57:LYS:N	30:BH:57:LYS:HE3	2.32	0.44
52:BI:52:LEU:HD12	52:BI:52:LEU:N	2.32	0.44
32:BK:121:GLU:HG2	37:BP:67:GLU:OE2	2.16	0.44
33:BL:58:TYR:O	33:BL:61:LEU:HD21	2.17	0.44
36:BO:8:ILE:O	36:BO:10:ARG:N	2.50	0.44
39:BR:35:PHE:HB2	39:BR:64:VAL:HG21	1.98	0.44
39:BR:47:VAL:C	39:BR:49:ILE:H	2.20	0.44
43:BW:28:GLU:C	43:BW:30:VAL:H	2.21	0.44
43:BW:39:GLN:HA	43:BW:68:PHE:HA	1.98	0.44
23:BB:1157:G:H1'	45:BY:10:ARG:NH2	2.31	0.44
1:CA:1030:U:HO2'	1:CA:1031:C:P	2.38	0.44
1:CA:1031:C:H5'	1:CA:1032:G:C4	2.53	0.44
1:CA:983:A:O2'	1:CA:1049:U:O2'	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:428:G:H4'	1:CA:429:U:O5'	2.17	0.44
1:CA:676:A:H5''	10:CK:114:PRO:CB	2.47	0.44
1:CA:836:G:H2'	1:CA:837:U:C6	2.52	0.44
1:CA:857:C:H2'	1:CA:858:G:O4'	2.17	0.44
2:CC:129:PHE:CZ	2:CC:156:LEU:HD13	2.52	0.44
4:CE:109:ALA:HB3	4:CE:135:VAL:HG11	1.99	0.44
4:CE:37:VAL:HG12	4:CE:116:VAL:HG21	1.99	0.44
4:CE:10:LEU:HA	4:CE:39:GLY:O	2.17	0.44
7:CH:76:ARG:NE	7:CH:125:ILE:HG23	2.32	0.44
7:CH:72:GLU:HB3	7:CH:129:ALA:HB2	1.98	0.44
8:CI:38:PHE:HB2	8:CI:44:ARG:HG3	2.00	0.44
9:CJ:17:LEU:HD11	9:CJ:96:VAL:HG12	1.98	0.44
9:CJ:98:VAL:O	9:CJ:98:VAL:HG23	2.18	0.44
12:CM:96:VAL:C	12:CM:98:GLY:H	2.20	0.44
14:CO:23:SER:HB3	14:CO:26:VAL:HG23	1.99	0.44
16:CQ:18:LYS:HA	16:CQ:47:ASP:O	2.17	0.44
48:D1:42:VAL:HG12	48:D1:43:ARG:HG2	1.99	0.44
22:DA:14:U:H4'	22:DA:70:C:O2	2.17	0.44
23:DB:152:A:H2'	23:DB:153:U:C6	2.53	0.44
23:DB:1563:U:O2'	23:DB:1564:C:H5'	2.17	0.44
23:DB:1724:G:H2'	23:DB:1725:U:H6	1.82	0.44
23:DB:2038:G:H2'	23:DB:2039:U:H6	1.82	0.44
23:DB:689:A:H2'	23:DB:690:G:C8	2.52	0.44
23:DB:709:U:H2'	23:DB:710:U:H6	1.81	0.44
23:DB:727:A:OP1	23:DB:1431:A:O2'	2.35	0.44
23:DB:807:U:H2'	23:DB:808:G:H8	1.81	0.44
23:DB:822:G:H2'	23:DB:823:C:C6	2.53	0.44
25:DC:137:GLY:O	25:DC:140:VAL:HG13	2.16	0.44
25:DC:155:ARG:NH2	25:DC:155:ARG:HG2	2.31	0.44
25:DC:24:HIS:HB3	25:DC:25:LYS:H	1.54	0.44
26:DD:202:ILE:HG22	26:DD:203:VAL:N	2.31	0.44
27:DE:109:LEU:HA	27:DE:117:ARG:NH1	2.32	0.44
27:DE:46:GLN:CD	27:DE:86:ALA:HB3	2.38	0.44
29:DG:98:LYS:HB3	29:DG:101:VAL:HG23	1.98	0.44
30:DH:35:LYS:NZ	30:DH:37:VAL:HG23	2.32	0.44
33:DL:58:TYR:HA	33:DL:62:PRO:CG	2.47	0.44
23:DB:2415:G:H4'	33:DL:66:PHE:HB2	2.00	0.44
34:DM:96:ILE:H	34:DM:98:PRO:HD3	1.83	0.44
38:DQ:92:LYS:C	38:DQ:93:ILE:HG23	2.36	0.44
39:DR:73:LYS:HD2	39:DR:73:LYS:N	2.32	0.44
45:DY:8:GLN:HE21	45:DY:23:LEU:HD12	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DZ:48:GLN:O	46:DZ:51:VAL:HB	2.17	0.44
1:AA:1085:U:H3'	1:AA:1086:U:C5	2.53	0.44
1:AA:1190:G:OP1	2:AC:4:VAL:HG12	2.17	0.44
1:AA:28:A:H2'	1:AA:29:U:O4'	2.17	0.44
1:AA:397:A:N7	1:AA:547:A:O2'	2.45	0.44
1:AA:541:G:H2'	1:AA:542:G:C8	2.52	0.44
20:AB:172:ILE:O	20:AB:175:ALA:HB3	2.17	0.44
2:AC:110:LEU:HB3	2:AC:203:LYS:NZ	2.32	0.44
2:AC:70:ALA:CB	2:AC:108:PRO:HB3	2.47	0.44
8:AI:118:ARG:HH22	8:AI:122:ARG:HH21	1.66	0.44
8:AI:51:LEU:CD1	8:AI:82:ILE:HG21	2.47	0.44
13:AN:73:LEU:HD12	13:AN:73:LEU:N	2.32	0.44
23:BB:1560:G:C4	23:BB:1561:C:C5	3.06	0.44
23:BB:1668:A:N3	23:BB:1670:C:C4	2.85	0.44
23:BB:1678:A:O2'	23:BB:1679:A:H5'	2.16	0.44
23:BB:1710:G:H2'	23:BB:1711:A:C8	2.52	0.44
23:BB:185:G:H2'	23:BB:186:G:O4'	2.17	0.44
23:BB:2100:G:H3'	23:BB:2101:A:H8	1.82	0.44
23:BB:2249:U:C4'	23:BB:2275:C:H5	2.29	0.44
23:BB:320:A:OP2	27:BE:132:LYS:HE2	2.18	0.44
23:BB:376:G:H2'	23:BB:377:G:H8	1.81	0.44
23:BB:39:G:H2'	23:BB:40:U:C6	2.52	0.44
23:BB:435:C:C2'	23:BB:436:C:H5'	2.47	0.44
23:BB:453:A:H4'	23:BB:472:A:N6	2.32	0.44
23:BB:608:A:H2'	23:BB:609:A:H8	1.81	0.44
23:BB:635:C:H3'	33:BL:109:LYS:NZ	2.33	0.44
23:BB:728:G:H5'	25:BC:12:ARG:HH22	1.82	0.44
25:BC:243:PRO:HB3	25:BC:249:VAL:HG23	1.99	0.44
23:BB:1802:A:O3'	25:BC:255:LYS:HE2	2.18	0.44
25:BC:7:PRO:HB3	25:BC:25:LYS:HD3	1.98	0.44
26:BD:13:ARG:O	26:BD:14:ILE:C	2.55	0.44
26:BD:62:LYS:H	26:BD:62:LYS:CE	2.30	0.44
29:BG:40:VAL:HG13	29:BG:52:GLY:O	2.18	0.44
31:BJ:119:PHE:O	31:BJ:122:LEU:HG	2.18	0.44
36:BO:10:ARG:NH1	36:BO:10:ARG:HG3	2.30	0.44
41:BT:47:VAL:HG12	41:BT:53:VAL:CG2	2.47	0.44
41:BT:62:VAL:O	41:BT:63:VAL:CG1	2.61	0.44
41:BT:87:LEU:HD22	41:BT:87:LEU:N	2.32	0.44
42:BU:42:LYS:HD3	42:BU:60:LYS:CB	2.45	0.44
24:BV:79:ARG:HB3	24:BV:79:ARG:NH1	2.32	0.44
43:BW:44:PHE:O	43:BW:76:ARG:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:BX:15:ASN:O	44:BX:17:GLU:N	2.51	0.44
45:BY:28:LEU:HA	45:BY:33:HIS:ND1	2.33	0.44
45:BY:9:THR:HB	45:BY:54:VAL:CA	2.39	0.44
1:CA:113:G:H2'	1:CA:114:U:H6	1.81	0.44
1:CA:1456:A:H2'	1:CA:1457:G:C8	2.52	0.44
1:CA:238:A:C3'	1:CA:239:U:H5''	2.47	0.44
1:CA:237:G:O2'	1:CA:238:A:H5'	2.18	0.44
1:CA:28:A:H2'	1:CA:29:U:O4'	2.17	0.44
1:CA:65:A:H2	1:CA:381:C:H2'	1.82	0.44
1:CA:598:U:H2'	1:CA:599:C:H6	1.82	0.44
1:CA:734:G:H2'	1:CA:735:C:C6	2.52	0.44
4:CE:64:GLU:HA	4:CE:67:ARG:NH1	2.32	0.44
5:CF:68:GLN:HA	5:CF:71:ILE:CG2	2.45	0.44
7:CH:75:GLN:HE21	7:CH:76:ARG:H	1.65	0.44
8:CI:30:ASN:O	8:CI:31:GLN:HB2	2.17	0.44
9:CJ:35:GLN:HB3	9:CJ:78:GLU:CD	2.37	0.44
10:CK:16:SER:N	10:CK:78:ILE:HA	2.32	0.44
12:CM:97:ARG:HG2	12:CM:97:ARG:HH11	1.82	0.44
13:CN:63:CYS:CB	13:CN:68:ARG:H	2.25	0.44
21:CU:43:GLU:O	21:CU:46:ARG:HB2	2.16	0.44
23:DB:1266:G:OP2	47:D0:12:ARG:NH1	2.50	0.44
51:D4:14:CYS:SG	51:D4:26:ILE:HG23	2.57	0.44
23:DB:1064:C:H2'	23:DB:1065:U:O4'	2.17	0.44
23:DB:1317:G:H2'	23:DB:1318:U:O4'	2.17	0.44
23:DB:1359:A:H2'	23:DB:1360:G:O4'	2.17	0.44
23:DB:1541:C:O2'	23:DB:1542:U:H5'	2.17	0.44
23:DB:1560:G:C4	23:DB:1561:C:C5	3.06	0.44
23:DB:1710:G:H2'	23:DB:1711:A:C8	2.52	0.44
23:DB:1964:G:H4'	23:DB:1965:C:OP2	2.17	0.44
23:DB:2270:A:H4'	43:DW:18:LYS:HB2	1.99	0.44
23:DB:2259:U:O4'	23:DB:2427:C:H2'	2.18	0.44
23:DB:2491:U:H5''	23:DB:2570:G:H5'	1.99	0.44
23:DB:458:G:O2'	23:DB:459:U:C5	2.69	0.44
23:DB:927:A:O2'	23:DB:928:A:H5'	2.18	0.44
25:DC:70:LYS:HD2	25:DC:95:TYR:CD1	2.52	0.44
26:DD:70:LYS:NZ	26:DD:70:LYS:HB3	2.32	0.44
29:DG:163:TYR:C	29:DG:165:ASP:H	2.20	0.44
31:DJ:93:ILE:HG22	31:DJ:97:PRO:HA	1.99	0.44
32:DK:17:ARG:HB2	32:DK:45:GLU:CB	2.47	0.44
32:DK:8:LEU:H	32:DK:8:LEU:HD12	1.82	0.44
36:DO:39:VAL:HG12	36:DO:50:ALA:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:47:ILE:HG23	37:DP:63:ILE:CA	2.32	0.44
23:DB:974:G:H4'	39:DR:83:TYR:CZ	2.52	0.44
41:DT:14:PRO:HA	41:DT:32:LEU:HG	1.98	0.44
42:DU:94:PHE:CD2	42:DU:95:PHE:N	2.84	0.44
24:DV:30:ILE:HB	24:DV:38:LEU:HB3	2.00	0.44
43:DW:19:ARG:NH2	43:DW:22:VAL:HA	2.32	0.44
1:AA:113:G:H2'	1:AA:114:U:H6	1.81	0.44
1:AA:1215:G:H5'	1:AA:1215:G:H8	1.80	0.44
1:AA:1296:C:O2'	1:AA:1302:C:C2	2.71	0.44
1:AA:1484:C:O2'	1:AA:1485:U:H5'	2.18	0.44
1:AA:397:A:N3	1:AA:397:A:H3'	2.33	0.44
1:AA:56:U:H6	1:AA:56:U:O5'	2.00	0.44
1:AA:60:A:O2'	1:AA:61:G:OP2	2.36	0.44
1:AA:951:G:H1'	1:AA:970:C:O2'	2.17	0.44
20:AB:162:VAL:CG1	20:AB:184:ALA:HB2	2.47	0.44
3:AD:120:LYS:CB	3:AD:145:ARG:HH21	2.30	0.44
10:AK:92:ARG:NH1	21:AU:24:LYS:NZ	2.65	0.44
11:AL:58:ASN:ND2	11:AL:60:PHE:HB2	2.33	0.44
12:AM:75:SER:O	12:AM:78:ARG:HB3	2.17	0.44
12:AM:78:ARG:HE	12:AM:79:LEU:CD1	2.29	0.44
16:AQ:22:VAL:HG21	16:AQ:58:VAL:HG21	1.99	0.44
17:AR:40:PRO:HD2	17:AR:43:ILE:HD11	1.98	0.44
18:AS:16:LYS:O	18:AS:20:LYS:HB2	2.17	0.44
23:BB:2393:U:OP1	50:B3:29:ARG:CZ	2.66	0.44
23:BB:1164:C:H2'	23:BB:1165:A:C8	2.52	0.44
23:BB:1189:A:H2'	23:BB:1190:G:O4'	2.17	0.44
23:BB:1210:G:H4'	23:BB:1211:C:OP2	2.18	0.44
23:BB:1228:G:H2'	23:BB:1229:C:H6	1.82	0.44
23:BB:144:A:H2'	23:BB:145:C:O4'	2.18	0.44
23:BB:121:G:H4'	23:BB:149:A:H5'	2.00	0.44
23:BB:1547:C:H2'	23:BB:1548:A:H8	1.81	0.44
23:BB:1559:U:H3'	23:BB:1560:G:C5'	2.47	0.44
23:BB:1346:G:C6	23:BB:1601:G:C6	3.05	0.44
23:BB:1785:A:H2'	23:BB:1787:A:C8	2.52	0.44
23:BB:966:G:H5''	23:BB:2272:U:O2	2.17	0.44
23:BB:2604:U:O2'	23:BB:2605:U:H5'	2.17	0.44
23:BB:68:G:N3	23:BB:68:G:H2'	2.31	0.44
23:BB:764:A:H5''	25:BC:208:GLY:HA3	1.98	0.44
23:BB:770:G:OP1	49:B2:8:SER:HB2	2.17	0.44
23:BB:863:A:H2'	23:BB:864:G:H8	1.83	0.44
23:BB:956:G:H22	23:BB:959:A:H3'	1.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:103:ILE:CG2	25:BC:104:LEU:N	2.80	0.44
26:BD:106:LYS:HB2	26:BD:106:LYS:HE3	1.77	0.44
26:BD:23:PRO:HA	26:BD:190:LYS:CA	2.47	0.44
28:BF:177:ARG:C	28:BF:177:ARG:HE	2.21	0.44
28:BF:77:LYS:HZ3	28:BF:79:ARG:HE	1.65	0.44
29:BG:24:THR:OG1	29:BG:31:GLU:HB3	2.17	0.44
29:BG:16:VAL:HG21	29:BG:44:HIS:NE2	2.32	0.44
30:BH:3:VAL:CG2	30:BH:21:VAL:HG11	2.44	0.44
52:BI:7:TYR:CD1	52:BI:7:TYR:C	2.91	0.44
32:BK:35:VAL:HG23	32:BK:36:GLY:N	2.22	0.44
35:BN:87:PHE:HB2	35:BN:94:TYR:CZ	2.52	0.44
22:BA:49:C:OP2	36:BO:30:ARG:NH2	2.50	0.44
37:BP:1:SER:HB2	37:BP:4:ILE:HB	1.98	0.44
39:BR:69:GLY:O	39:BR:70:GLU:HG3	2.17	0.44
39:BR:78:ARG:HD3	39:BR:87:GLN:HG2	2.00	0.44
40:BS:50:VAL:HG21	40:BS:103:ILE:HG21	1.98	0.44
43:BW:48:ALA:HA	43:BW:54:ARG:HA	2.00	0.44
43:BW:49:ASN:O	43:BW:50:VAL:HB	2.18	0.44
1:CA:11:G:H2'	1:CA:12:U:C6	2.53	0.44
1:CA:178:C:O2'	1:CA:179:A:H5'	2.18	0.44
1:CA:29:U:O2'	1:CA:30:U:H5'	2.17	0.44
1:CA:36:C:H2'	1:CA:37:U:O4'	2.17	0.44
1:CA:668:G:O2'	1:CA:669:G:H5'	2.17	0.44
1:CA:81:A:H2'	1:CA:82:G:H8	1.81	0.44
2:CC:108:PRO:C	2:CC:110:LEU:H	2.21	0.44
2:CC:65:VAL:O	2:CC:65:VAL:HG12	2.17	0.44
2:CC:87:ARG:HG3	2:CC:87:ARG:HH11	1.81	0.44
3:CD:102:TYR:HE1	3:CD:108:ALA:O	2.01	0.44
3:CD:55:ARG:HH11	3:CD:55:ARG:CG	2.31	0.44
12:CM:21:ILE:O	12:CM:21:ILE:HG22	2.16	0.44
47:D0:47:TYR:CD2	47:D0:51:ARG:HA	2.53	0.44
51:D4:30:GLU:O	51:D4:32:LYS:N	2.50	0.44
22:DA:49:C:H2'	22:DA:50:A:H8	1.83	0.44
23:DB:1128:G:N7	23:DB:2490:G:H5'	2.32	0.44
23:DB:1292:G:H2'	23:DB:1293:C:H6	1.83	0.44
23:DB:145:C:H2'	23:DB:146:A:C8	2.52	0.44
23:DB:20:C:H2'	23:DB:21:A:C8	2.51	0.44
23:DB:2215:C:O2'	23:DB:2216:G:H5'	2.17	0.44
23:DB:858:G:C2	23:DB:2268:A:H2'	2.47	0.44
23:DB:2377:A:H2	36:DO:92:PHE:HE1	1.65	0.44
23:DB:239:C:H2'	23:DB:240:C:C6	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2521:C:O2'	23:DB:2522:U:H5'	2.17	0.44
23:DB:2669:G:H2'	23:DB:2670:A:C8	2.51	0.44
23:DB:2758:A:H2'	23:DB:2759:G:C5'	2.45	0.44
23:DB:304:U:H2'	23:DB:305:C:H6	1.78	0.44
23:DB:357:C:H2'	23:DB:358:U:H6	1.82	0.44
23:DB:383:C:H5''	23:DB:385:C:OP2	2.17	0.44
23:DB:182:A:H2	23:DB:433:C:O2	2.00	0.44
23:DB:538:A:N6	23:DB:555:G:O2'	2.43	0.44
23:DB:7:G:H2'	23:DB:8:C:H6	1.82	0.44
23:DB:814:C:H5''	39:DR:87:GLN:O	2.18	0.44
22:DA:99:A:N1	23:DB:863:A:H4'	2.33	0.44
26:DD:122:VAL:CG2	26:DD:128:ARG:HG3	2.46	0.44
26:DD:83:ARG:HD2	26:DD:84:LEU:N	2.32	0.44
27:DE:148:ILE:HG13	27:DE:148:ILE:O	2.17	0.44
27:DE:151:GLY:HA2	27:DE:169:VAL:O	2.17	0.44
29:DG:156:TYR:CD1	29:DG:171:LYS:HG2	2.53	0.44
33:DL:61:LEU:HD23	33:DL:61:LEU:C	2.38	0.44
34:DM:33:LEU:CD2	34:DM:124:LEU:HD22	2.48	0.44
37:DP:8:GLU:HA	37:DP:11:GLN:HG2	1.99	0.44
40:DS:18:ARG:NE	40:DS:22:ASP:OD1	2.50	0.44
40:DS:22:ASP:O	40:DS:25:ARG:HG3	2.18	0.44
24:DV:72:VAL:CG1	24:DV:93:ARG:HA	2.42	0.44
1:AA:1148:U:H5''	8:AI:8:THR:HG23	2.00	0.44
1:AA:1521:C:H2'	1:AA:1522:U:C6	2.53	0.44
1:AA:1529:G:OP2	1:AA:1529:G:H3'	2.18	0.44
1:AA:613:C:H2'	1:AA:614:C:H6	1.82	0.44
1:AA:915:A:C2'	1:AA:916:U:H5'	2.48	0.44
20:AB:9:LEU:H	20:AB:9:LEU:HD12	1.82	0.44
2:AC:148:ILE:HA	2:AC:200:TRP:O	2.18	0.44
3:AD:190:LEU:O	3:AD:192:ALA:N	2.50	0.44
3:AD:12:ARG:HD2	3:AD:37:PRO:HA	2.00	0.44
4:AE:105:ILE:HD12	4:AE:123:LEU:HB3	2.00	0.44
8:AI:4:GLN:NE2	8:AI:21:LYS:HE3	2.33	0.44
9:AJ:57:VAL:HG22	9:AJ:58:ASN:H	1.83	0.44
13:AN:17:ASP:HB2	13:AN:18:LYS:NZ	2.33	0.44
13:AN:49:THR:O	13:AN:49:THR:HG22	2.17	0.44
13:AN:63:CYS:HB2	13:AN:79:SER:OG	2.18	0.44
15:AP:78:VAL:O	15:AP:80:LYS:N	2.51	0.44
18:AS:35:ARG:HG3	18:AS:35:ARG:H	1.50	0.44
18:AS:47:THR:HG23	18:AS:60:PHE:HE1	1.83	0.44
23:BB:1107:G:H2'	23:BB:1108:U:C6	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1231:U:H2'	23:BB:1232:G:C8	2.52	0.44
23:BB:1535:A:H5'	23:BB:1536:C:C5	2.53	0.44
23:BB:1561:C:H2'	23:BB:1562:U:H6	1.81	0.44
23:BB:1311:G:H21	23:BB:1603:A:H62	1.65	0.44
23:BB:1877:A:O2'	23:BB:1878:G:H5'	2.18	0.44
23:BB:1910:G:O2'	23:BB:1911:U:H5'	2.18	0.44
23:BB:1964:G:H4'	23:BB:1965:C:OP2	2.17	0.44
23:BB:2150:C:H2'	23:BB:2151:U:O4'	2.17	0.44
23:BB:230:G:H2'	23:BB:231:A:H8	1.82	0.44
23:BB:1027:A:H2	23:BB:2488:G:H4'	1.82	0.44
23:BB:2669:G:H2'	23:BB:2670:A:C8	2.51	0.44
23:BB:2787:C:H2'	23:BB:2788:C:C6	2.53	0.44
23:BB:661:A:H2'	23:BB:662:G:H8	1.83	0.44
23:BB:743:A:C2'	23:BB:744:U:H5'	2.47	0.44
23:BB:985:C:H2'	23:BB:986:C:H6	1.83	0.44
25:BC:124:LYS:HZ3	25:BC:125:PRO:HD2	1.82	0.44
25:BC:216:ARG:HB3	25:BC:217:PRO:HD2	1.99	0.44
25:BC:6:LYS:HA	25:BC:7:PRO:HA	1.82	0.44
26:BD:171:THR:HG22	26:BD:173:GLN:HG3	1.98	0.44
26:BD:29:VAL:HA	26:BD:185:ASN:OD1	2.18	0.44
26:BD:40:LEU:O	26:BD:41:ALA:HB3	2.18	0.44
27:BE:111:GLU:C	27:BE:113:VAL:H	2.21	0.44
27:BE:178:VAL:HG23	27:BE:179:SER:N	2.33	0.44
28:BF:116:LEU:H	28:BF:176:PHE:CA	2.31	0.44
28:BF:48:LEU:O	28:BF:49:LEU:HB2	2.17	0.44
31:BJ:101:ILE:HG12	31:BJ:102:GLU:N	2.32	0.44
23:BB:1669:A:C8	32:BK:5:GLN:HG3	2.53	0.44
32:BK:75:SER:N	37:BP:76:HIS:NE2	2.64	0.44
33:BL:78:ARG:NH1	33:BL:80:SER:H	2.16	0.44
35:BN:24:MET:HE2	35:BN:44:LEU:HD13	1.99	0.44
40:BS:70:LYS:CD	40:BS:110:ARG:HG3	2.47	0.44
40:BS:9:HIS:O	40:BS:10:ALA:HB3	2.18	0.44
42:BU:2:ALA:O	42:BU:4:ILE:HG13	2.18	0.44
24:BV:77:VAL:HG12	34:BM:133:LYS:HZ2	1.83	0.44
43:BW:37:VAL:O	43:BW:38:ARG:C	2.55	0.44
1:CA:1142:G:H2'	1:CA:1143:G:O4'	2.17	0.44
1:CA:1300:G:H2'	1:CA:1301:U:OP2	2.18	0.44
1:CA:1349:A:H61	1:CA:1373:G:H1'	1.82	0.44
1:CA:1414:U:H2'	1:CA:1415:G:C8	2.52	0.44
1:CA:1425:U:H2'	1:CA:1426:G:C8	2.53	0.44
1:CA:432:A:H2'	1:CA:433:G:H5'	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:949:A:O2'	1:CA:950:U:H5'	2.18	0.44
2:CC:53:ARG:HA	2:CC:113:LYS:HZ3	1.83	0.44
3:CD:107:GLY:O	3:CD:157:ALA:HB1	2.17	0.44
3:CD:37:PRO:HD2	3:CD:41:GLY:CA	2.47	0.44
5:CF:72:ASP:O	5:CF:75:GLU:HB3	2.18	0.44
8:CI:6:TYR:OH	8:CI:8:THR:HG22	2.17	0.44
9:CJ:68:ARG:HD3	9:CJ:70:HIS:CE1	2.53	0.44
10:CK:117:HIS:O	10:CK:118:ASN:HB2	2.16	0.44
10:CK:12:ARG:CB	10:CK:13:LYS:HD2	2.47	0.44
10:CK:70:ALA:C	10:CK:72:ALA:N	2.70	0.44
10:CK:84:MET:SD	10:CK:110:THR:OG1	2.73	0.44
11:CL:19:ASN:HB3	11:CL:85:ARG:HD2	1.99	0.44
14:CO:42:PHE:CD1	14:CO:55:LEU:HD22	2.53	0.44
16:CQ:68:LYS:C	16:CQ:70:LYS:N	2.70	0.44
17:CR:42:ARG:C	17:CR:44:THR:H	2.21	0.44
18:CS:10:ILE:CG2	18:CS:40:PHE:HE2	2.31	0.44
47:D0:36:LYS:HG3	47:D0:40:HIS:C	2.38	0.44
51:D4:23:ILE:HD12	51:D4:24:ARG:HH11	1.82	0.44
23:DB:1176:U:H2'	23:DB:1177:G:O4'	2.17	0.44
23:DB:11:C:H2'	23:DB:12:U:C5'	2.44	0.44
23:DB:136:G:N2	41:DT:3:ARG:CZ	2.80	0.44
23:DB:1341:G:C2'	23:DB:1397:U:O2'	2.66	0.44
23:DB:1541:C:H2'	23:DB:1542:U:O4'	2.16	0.44
23:DB:1589:U:H2'	23:DB:1590:A:C8	2.52	0.44
23:DB:1659:G:C6	23:DB:2002:G:C6	3.06	0.44
23:DB:1775:U:H2'	23:DB:1776:G:C5'	2.47	0.44
23:DB:1785:A:H2'	23:DB:1787:A:C8	2.52	0.44
23:DB:1930:G:HO2'	23:DB:1931:U:H5	1.64	0.44
23:DB:198:C:H5'	23:DB:2244:U:OP1	2.18	0.44
23:DB:2294:G:OP2	36:DO:9:ARG:NE	2.49	0.44
23:DB:2321:U:O2	23:DB:2321:U:O4'	2.35	0.44
23:DB:2463:C:O2'	23:DB:2464:G:H5'	2.18	0.44
23:DB:2533:U:H2'	23:DB:2534:A:O4'	2.17	0.44
23:DB:2728:U:H2'	23:DB:2729:G:H8	1.82	0.44
23:DB:2867:G:N3	23:DB:2867:G:C2'	2.73	0.44
23:DB:2887:A:O4'	47:D0:27:LEU:HD11	2.17	0.44
23:DB:298:G:N2	23:DB:339:U:C5	2.86	0.44
23:DB:319:G:H2'	23:DB:320:A:O4'	2.18	0.44
23:DB:401:A:O2'	23:DB:402:A:C5'	2.65	0.44
25:DC:135:PRO:O	25:DC:165:ALA:HA	2.17	0.44
25:DC:23:LEU:HD12	25:DC:24:HIS:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DG:93:TYR:HA	29:DG:105:SER:O	2.17	0.44
29:DG:134:GLY:N	29:DG:140:ILE:HD11	2.32	0.44
30:DH:108:VAL:HG12	30:DH:110:VAL:CB	2.47	0.44
30:DH:4:ILE:CD1	30:DH:4:ILE:H	2.31	0.44
33:DL:64:PHE:CD2	50:D3:11:LYS:HB3	2.49	0.44
34:DM:119:LEU:N	34:DM:119:LEU:HD22	2.32	0.44
34:DM:5:LYS:HE3	34:DM:6:ARG:CA	2.45	0.44
35:DN:121:LYS:HD2	35:DN:122:ALA:H	1.83	0.44
37:DP:34:GLY:O	37:DP:35:SER:HB3	2.17	0.44
38:DQ:32:ARG:NH1	38:DQ:33:VAL:HG22	2.33	0.44
40:DS:42:LYS:HE2	40:DS:45:VAL:HG21	1.99	0.44
41:DT:9:LYS:HD3	44:DX:29:ARG:HH22	1.82	0.44
46:DZ:2:LYS:O	46:DZ:3:LYS:C	2.55	0.44
1:AA:1340:A:H2'	1:AA:1341:U:H6	1.81	0.44
1:AA:1347:G:O2'	1:AA:1348:U:P	2.75	0.44
1:AA:206:C:H2'	1:AA:207:C:C6	2.53	0.44
1:AA:500:G:H2'	1:AA:501:C:C6	2.53	0.44
20:AB:104:LYS:HE3	20:AB:104:LYS:HB2	1.82	0.44
2:AC:41:TYR:HA	2:AC:44:LYS:CD	2.47	0.44
2:AC:77:GLY:CA	2:AC:81:GLU:HB3	2.46	0.44
3:AD:39:GLN:HB3	3:AD:40:HIS:ND1	2.32	0.44
3:AD:71:PHE:O	3:AD:74:TYR:HB2	2.17	0.44
6:AG:115:MET:CE	6:AG:119:LEU:HB2	2.47	0.44
7:AH:102:VAL:HG12	7:AH:125:ILE:HD12	1.99	0.44
10:AK:109:ILE:HG22	21:AU:16:ARG:HD2	1.99	0.44
15:AP:6:LEU:HD12	15:AP:6:LEU:N	2.32	0.44
16:AQ:39:ARG:HH11	16:AQ:39:ARG:CB	2.30	0.44
23:BB:443:A:C2	23:BB:1245:G:N3	2.86	0.44
23:BB:137:U:C2'	23:BB:138:U:H5'	2.48	0.44
23:BB:1418:G:N2	23:BB:1579:A:C8	2.85	0.44
23:BB:1509:A:C4'	23:BB:1510:G:H5'	2.48	0.44
23:BB:1552:A:H2'	23:BB:1553:A:C5'	2.44	0.44
23:BB:1724:G:H2'	23:BB:1725:U:H6	1.82	0.44
23:BB:197:A:H62	23:BB:2430:A:H2'	1.83	0.44
23:BB:2206:C:O2'	23:BB:2207:C:H5'	2.17	0.44
23:BB:2236:U:O2'	23:BB:2237:G:H5'	2.18	0.44
23:BB:2508:G:O3'	23:BB:2555:U:H5'	2.18	0.44
23:BB:2712:C:H2'	23:BB:2714:G:O3'	2.17	0.44
23:BB:346:A:H3'	23:BB:347:A:H8	1.83	0.44
23:BB:378:C:O2'	23:BB:379:G:H5'	2.17	0.44
23:BB:43:G:H2'	23:BB:44:A:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:565:C:O2'	23:BB:566:U:H5'	2.18	0.44
25:BC:153:LEU:CD1	25:BC:161:VAL:HG23	2.47	0.44
25:BC:237:ARG:HB2	25:BC:239:PHE:CE1	2.52	0.44
23:BB:2512:C:O2'	26:BD:146:ILE:CG1	2.66	0.44
27:BE:148:ILE:N	27:BE:183:PHE:HB3	2.31	0.44
28:BF:107:VAL:HB	28:BF:108:PRO:HD3	2.00	0.44
29:BG:125:PRO:HD2	29:BG:129:GLU:HB2	2.00	0.44
29:BG:44:HIS:HB2	29:BG:49:LEU:HD23	2.00	0.44
30:BH:23:ALA:C	30:BH:25:TYR:H	2.21	0.44
31:BJ:71:ASP:O	31:BJ:72:LYS:HB2	2.17	0.44
31:BJ:85:LYS:HD3	31:BJ:85:LYS:HA	1.77	0.44
34:BM:45:GLN:HE21	34:BM:119:LEU:HB3	1.81	0.44
34:BM:58:LYS:C	34:BM:60:GLN:N	2.71	0.44
35:BN:9:GLN:HB2	35:BN:9:GLN:HE21	1.54	0.44
36:BO:76:LYS:HB2	36:BO:106:LEU:HD13	1.99	0.44
37:BP:18:SER:C	37:BP:20:ARG:H	2.21	0.44
37:BP:18:SER:OG	37:BP:21:PRO:HD3	2.18	0.44
39:BR:42:ALA:HA	39:BR:54:VAL:HG22	1.99	0.44
39:BR:68:ARG:HD2	39:BR:70:GLU:OE1	2.17	0.44
40:BS:17:VAL:HG23	40:BS:76:VAL:HG21	2.00	0.44
40:BS:86:MET:HB2	40:BS:96:ILE:HD11	1.98	0.44
41:BT:15:HIS:HB2	41:BT:31:VAL:HG21	1.99	0.44
41:BT:6:ARG:O	41:BT:8:LEU:N	2.50	0.44
41:BT:71:GLY:O	41:BT:72:GLN:HG3	2.17	0.44
23:BB:456:C:H3'	41:BT:72:GLN:HE22	1.82	0.44
42:BU:13:LEU:N	42:BU:68:ASN:ND2	2.64	0.44
42:BU:11:ILE:O	42:BU:69:VAL:HB	2.18	0.44
24:BV:30:ILE:O	24:BV:37:PRO:HA	2.18	0.44
43:BW:43:LYS:HA	43:BW:63:ASP:OD1	2.18	0.44
43:BW:66:VAL:CG2	43:BW:67:LYS:H	2.26	0.44
46:BZ:63:ARG:NH2	46:BZ:66:ILE:HD12	2.33	0.44
1:CA:1299:A:C5	1:CA:1301:U:H1'	2.53	0.44
1:CA:1481:U:H2'	1:CA:1482:G:C8	2.52	0.44
1:CA:234:C:H2'	1:CA:235:C:C6	2.53	0.44
1:CA:239:U:C6	1:CA:239:U:C5'	2.97	0.44
1:CA:448:A:H2'	1:CA:449:G:C8	2.53	0.44
1:CA:659:U:O2'	1:CA:660:C:H5'	2.18	0.44
1:CA:729:A:H2'	1:CA:730:G:H8	1.83	0.44
1:CA:731:G:OP1	1:CA:766:A:H1'	2.16	0.44
1:CA:86:G:O2'	1:CA:87:C:P	2.76	0.44
2:CC:156:LEU:CD1	2:CC:165:GLU:HB2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CD:100:VAL:HG12	3:CD:101:VAL:N	2.33	0.44
4:CE:99:SER:C	4:CE:101:GLY:H	2.20	0.44
6:CG:147:ASN:C	6:CG:149:ALA:H	2.21	0.44
6:CG:78:ARG:HD2	6:CG:81:GLY:H	1.82	0.44
7:CH:103:VAL:HG22	7:CH:124:ILE:HA	1.99	0.44
1:CA:705:G:N2	10:CK:43:TRP:CE3	2.85	0.44
1:CA:695:A:C5'	10:CK:52:ARG:HH22	2.29	0.44
15:CP:40:ASN:HD21	15:CP:42:ILE:CG1	2.30	0.44
18:CS:5:LYS:C	18:CS:6:LYS:HG2	2.38	0.44
22:DA:92:C:H2'	22:DA:93:C:H6	1.82	0.44
23:DB:1028:A:N6	23:DB:1125:G:H2'	2.33	0.44
23:DB:1145:C:O2'	23:DB:1146:C:H5'	2.17	0.44
23:DB:1529:G:H2'	23:DB:1530:G:C8	2.53	0.44
23:DB:1746:A:H2'	23:DB:1747:U:C6	2.53	0.44
23:DB:1985:C:O2'	23:DB:1986:C:H5'	2.17	0.44
23:DB:2438:U:O2'	23:DB:2439:A:H5''	2.18	0.44
23:DB:1027:A:N3	23:DB:2488:G:H5''	2.32	0.44
23:DB:435:C:H2'	23:DB:436:C:H5'	1.99	0.44
23:DB:635:C:OP2	33:DL:126:ARG:NE	2.42	0.44
23:DB:661:A:H2'	23:DB:662:G:H8	1.82	0.44
25:DC:51:ARG:HD3	25:DC:51:ARG:O	2.17	0.44
26:DD:180:VAL:HG23	26:DD:181:ASP:N	2.32	0.44
27:DE:147:LEU:O	27:DE:148:ILE:HG13	2.18	0.44
52:DI:138:VAL:HG12	52:DI:139:VAL:N	2.32	0.44
52:DI:54:ILE:C	52:DI:54:ILE:HD13	2.37	0.44
31:DJ:110:PRO:O	31:DJ:111:LYS:HB2	2.18	0.44
31:DJ:73:VAL:HG21	31:DJ:75:TYR:CE2	2.52	0.44
31:DJ:94:ALA:CB	31:DJ:95:ARG:HH21	2.31	0.44
33:DL:120:VAL:HG12	33:DL:122:VAL:HG22	1.98	0.44
36:DO:79:ALA:O	36:DO:82:ALA:HB3	2.18	0.44
40:DS:1:MET:HB3	40:DS:4:ILE:CD1	2.47	0.44
40:DS:43:ALA:C	40:DS:45:VAL:H	2.20	0.44
40:DS:42:LYS:O	40:DS:45:VAL:HG22	2.17	0.44
40:DS:84:ARG:HH21	40:DS:98:LYS:HZ1	1.66	0.44
44:DX:22:LEU:HD22	44:DX:25:GLN:CD	2.38	0.44
1:AA:1001:C:H2'	1:AA:1002:G:C8	2.53	0.44
1:AA:1050:G:O2'	1:AA:1051:C:H5'	2.18	0.44
1:AA:1151:A:HO2'	1:AA:1152:A:H8	1.61	0.44
1:AA:1264:U:O2'	1:AA:1265:C:H5'	2.17	0.44
1:AA:29:U:H5'	1:AA:296:U:OP1	2.17	0.44
1:AA:993:G:C2'	1:AA:995:C:H41	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:32:LYS:O	3:AD:35:GLN:HB2	2.17	0.44
5:AF:41:ASP:OD1	5:AF:58:HIS:HE1	2.01	0.44
5:AF:6:ILE:HB	5:AF:89:VAL:HB	2.00	0.44
8:AI:43:ALA:C	8:AI:45:MET:H	2.21	0.44
10:AK:69:CYS:C	10:AK:71:ASP:N	2.70	0.44
12:AM:44:ILE:O	12:AM:47:LEU:HB2	2.17	0.44
13:AN:11:LYS:O	13:AN:15:LEU:HG	2.18	0.44
9:AJ:66:GLU:OE2	13:AN:98:ALA:HA	2.17	0.44
17:AR:63:TYR:C	17:AR:65:SER:H	2.21	0.44
19:AT:27:MET:HE3	19:AT:66:ILE:HG12	1.99	0.44
23:BB:1110:G:N2	23:BB:1111:A:N6	2.66	0.44
23:BB:1585:C:H2'	23:BB:1586:A:C8	2.52	0.44
23:BB:1843:C:H2'	23:BB:1844:C:H6	1.81	0.44
23:BB:2055:C:H4'	23:BB:2056:G:H5''	1.99	0.44
23:BB:2070:A:H2'	23:BB:2071:A:O4'	2.18	0.44
23:BB:2221:G:H2'	23:BB:2222:C:H6	1.83	0.44
23:BB:2286:G:H5'	23:BB:2286:G:N3	2.33	0.44
23:BB:2425:A:O2'	23:BB:2426:A:OP2	2.28	0.44
23:BB:2601:C:C3'	23:BB:2602:A:H5''	2.48	0.44
23:BB:2617:U:H2'	23:BB:2618:G:C5'	2.48	0.44
23:BB:2733:A:H2'	23:BB:2734:A:O4'	2.17	0.44
23:BB:296:U:H2'	23:BB:297:G:C8	2.48	0.44
23:BB:455:C:N3	23:BB:472:A:H2'	2.33	0.44
23:BB:512:G:H4'	23:BB:512:G:OP1	2.18	0.44
25:BC:103:ILE:HG22	25:BC:104:LEU:H	1.81	0.44
25:BC:140:VAL:HG23	25:BC:193:GLU:OE2	2.17	0.44
26:BD:135:GLY:C	26:BD:137:SER:H	2.21	0.44
26:BD:2:ILE:CG2	26:BD:204:LYS:HG3	2.48	0.44
26:BD:88:GLU:HG2	26:BD:89:GLU:H	1.83	0.44
27:BE:12:LEU:CD2	27:BE:118:LEU:HA	2.48	0.44
23:BB:675:A:O2'	27:BE:62:GLN:NE2	2.51	0.44
52:BI:11:GLN:NE2	52:BI:74:PRO:HG2	2.33	0.44
52:BI:7:TYR:HB2	52:BI:59:THR:HA	1.99	0.44
32:BK:63:VAL:HG13	32:BK:107:LEU:HD22	2.00	0.44
34:BM:8:LYS:O	34:BM:10:ARG:N	2.50	0.44
35:BN:45:ARG:HD2	35:BN:45:ARG:H	1.82	0.44
36:BO:27:VAL:HA	36:BO:94:ARG:HB3	2.00	0.44
37:BP:24:THR:HG21	37:BP:111:GLU:OE2	2.17	0.44
37:BP:92:ARG:HB3	37:BP:110:LYS:O	2.18	0.44
39:BR:41:ILE:HG12	39:BR:60:LYS:O	2.17	0.44
24:BV:79:ARG:NH1	24:BV:86:LEU:HD11	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1172:C:O2'	1:CA:1173:U:H5'	2.17	0.44
1:CA:223:A:H2'	1:CA:224:U:C6	2.53	0.44
1:CA:109:A:C6	1:CA:326:G:C6	3.06	0.44
1:CA:397:A:H3'	1:CA:397:A:N3	2.33	0.44
1:CA:401:C:O2'	1:CA:402:G:H5'	2.18	0.44
1:CA:500:G:H2'	1:CA:501:C:C6	2.53	0.44
1:CA:542:G:O2'	1:CA:543:U:H5'	2.18	0.44
1:CA:605:U:H2'	1:CA:606:G:C8	2.53	0.44
1:CA:89:U:O2'	1:CA:90:C:H5'	2.17	0.44
20:CB:72:LYS:O	20:CB:74:ALA:N	2.51	0.44
3:CD:137:SER:HB3	3:CD:138:PRO:HD2	2.00	0.44
3:CD:169:TRP:CD1	3:CD:170:LEU:HG	2.53	0.44
5:CF:30:THR:HA	5:CF:34:GLY:O	2.18	0.44
8:CI:42:THR:HA	8:CI:45:MET:SD	2.57	0.44
8:CI:5:TYR:O	8:CI:19:PHE:HB3	2.16	0.44
1:CA:947:G:H5''	12:CM:106:ARG:HB2	2.00	0.44
12:CM:95:PRO:HG2	12:CM:101:THR:HG22	1.98	0.44
13:CN:71:GLY:O	13:CN:79:SER:HA	2.17	0.44
14:CO:70:LYS:HB2	14:CO:77:TYR:CG	2.52	0.44
1:CA:719:C:O2	17:CR:37:LYS:HA	2.18	0.44
18:CS:39:ILE:HD11	18:CS:70:LEU:HD13	2.00	0.44
47:D0:23:ALA:C	47:D0:25:THR:H	2.21	0.44
23:DB:104:A:H2'	23:DB:105:C:C6	2.49	0.44
23:DB:1098:A:C8	52:DI:3:LYS:HB3	2.53	0.44
23:DB:1171:G:N1	23:DB:1179:G:C6	2.85	0.44
23:DB:1446:C:H2'	23:DB:1447:C:C6	2.52	0.44
23:DB:165:A:H2'	23:DB:166:U:C6	2.50	0.44
23:DB:1680:U:H2'	23:DB:1681:G:O4'	2.18	0.44
23:DB:245:G:OP1	33:DL:67:THR:HB	2.18	0.44
23:DB:619:G:H2'	23:DB:620:G:H5''	2.00	0.44
23:DB:729:G:H4'	23:DB:763:G:C5'	2.47	0.44
25:DC:251:THR:O	25:DC:252:LYS:CB	2.65	0.44
23:DB:1802:A:C4'	25:DC:255:LYS:HE2	2.40	0.44
26:DD:24:VAL:HG21	26:DD:193:VAL:HG13	2.00	0.44
26:DD:27:ILE:C	26:DD:27:ILE:HD13	2.37	0.44
27:DE:58:LYS:HA	27:DE:59:PRO:HD3	1.66	0.44
28:DF:35:LEU:HA	28:DF:153:ILE:HA	2.00	0.44
29:DG:42:VAL:HG22	29:DG:43:LYS:N	2.33	0.44
30:DH:104:THR:HA	30:DH:108:VAL:O	2.18	0.44
30:DH:114:GLU:HB3	30:DH:133:GLN:NE2	2.25	0.44
52:DI:7:TYR:CZ	52:DI:57:VAL:HG11	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DJ:7:LYS:HD2	31:DJ:45:THR:HG21	1.99	0.44
31:DJ:4:PHE:HD1	31:DJ:5:THR:N	2.16	0.44
32:DK:2:ILE:HG23	32:DK:33:ALA:O	2.17	0.44
23:DB:1243:C:H1'	33:DL:14:LYS:HE3	2.00	0.44
38:DQ:88:GLU:O	38:DQ:89:ILE:C	2.56	0.44
40:DS:56:ALA:O	40:DS:59:GLU:HB3	2.17	0.44
40:DS:92:ARG:HH11	40:DS:92:ARG:HG2	1.82	0.44
40:DS:9:HIS:HA	40:DS:100:THR:OG1	2.18	0.44
41:DT:15:HIS:O	41:DT:16:VAL:CB	2.66	0.44
23:DB:144:A:C5	41:DT:3:ARG:NH1	2.86	0.44
42:DU:8:ASP:HB2	42:DU:25:LYS:HE3	2.00	0.44
24:DV:29:ILE:HG13	24:DV:30:ILE:N	2.33	0.44
43:DW:45:HIS:HE1	43:DW:71:LYS:HG3	1.81	0.44
43:DW:81:ILE:HG23	43:DW:83:ALA:H	1.82	0.44
45:DY:43:ILE:HA	45:DY:46:MET:CB	2.48	0.44
1:AA:1172:C:H2'	1:AA:1173:U:H6	1.82	0.44
1:AA:11:G:H2'	1:AA:12:U:C6	2.52	0.44
1:AA:15:G:O2'	4:AE:21:SER:HB2	2.17	0.44
1:AA:432:A:H2'	1:AA:433:G:H5'	1.99	0.44
1:AA:488:C:H2'	1:AA:489:C:C6	2.52	0.44
1:AA:857:C:H2'	1:AA:858:G:O4'	2.17	0.44
1:AA:989:U:O2'	1:AA:990:C:H5'	2.18	0.44
20:AB:131:LYS:O	20:AB:134:LEU:HB2	2.18	0.44
20:AB:90:PHE:CE2	20:AB:149:GLY:HA3	2.52	0.44
2:AC:171:ARG:HH12	2:AC:173:PRO:HG3	1.83	0.44
5:AF:38:ARG:CB	5:AF:63:ASN:HB2	2.37	0.44
8:AI:33:SER:HB3	8:AI:36:GLN:CB	2.43	0.44
8:AI:43:ALA:HA	8:AI:44:ARG:HH12	1.82	0.44
9:AJ:37:ARG:HB2	9:AJ:77:VAL:CG1	2.48	0.44
9:AJ:36:VAL:HG22	9:AJ:76:ILE:HG22	1.99	0.44
9:AJ:11:LYS:NZ	9:AJ:99:GLN:H	2.15	0.44
11:AL:117:GLY:O	11:AL:118:VAL:HG13	2.18	0.44
11:AL:49:ARG:C	11:AL:50:LYS:HE2	2.38	0.44
1:AA:1049:U:C2'	13:AN:2:LYS:HD3	2.45	0.44
18:AS:44:ILE:HA	18:AS:61:VAL:CG1	2.48	0.44
47:B0:35:GLU:C	47:B0:37:HIS:H	2.21	0.44
48:B1:10:LEU:CD1	48:B1:25:ASN:HD22	2.31	0.44
23:BB:55:G:P	49:B2:35:ARG:HE	2.40	0.44
22:BA:45:A:O2'	22:BA:46:A:H5'	2.18	0.44
22:BA:64:G:H2'	22:BA:65:U:C6	2.53	0.44
23:BB:1080:A:H2'	23:BB:1081:U:C6	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1152:C:H3'	55:BB:3244:HOH:O	2.17	0.44
23:BB:1529:G:H2'	23:BB:1530:G:C8	2.52	0.44
23:BB:1592:C:H2'	23:BB:1593:A:H8	1.83	0.44
23:BB:1965:C:H3'	23:BB:1966:A:H8	1.83	0.44
23:BB:2247:A:H3'	55:BB:3268:HOH:O	2.17	0.44
23:BB:2305:U:N3	28:BF:150:GLY:HA3	2.33	0.44
23:BB:2324:U:C5'	23:BB:2325:G:H5''	2.45	0.44
23:BB:2581:G:N3	23:BB:2581:G:H2'	2.33	0.44
23:BB:338:G:C2'	23:BB:339:U:H5'	2.48	0.44
23:BB:340:A:C2'	23:BB:341:C:H5'	2.48	0.44
25:BC:128:THR:HG22	25:BC:129:LEU:H	1.83	0.44
25:BC:198:GLU:C	25:BC:200:MET:N	2.70	0.44
26:BD:155:VAL:O	26:BD:156:PHE:HB2	2.18	0.44
26:BD:31:ALA:O	26:BD:32:ASN:ND2	2.50	0.44
27:BE:122:GLU:OE2	27:BE:152:GLU:HB2	2.18	0.44
27:BE:117:ARG:HA	27:BE:188:MET:SD	2.57	0.44
27:BE:40:ARG:HG2	27:BE:92:HIS:HB2	2.00	0.44
30:BH:30:LEU:O	30:BH:35:LYS:N	2.50	0.44
32:BK:26:GLY:HA3	32:BK:30:ARG:HG3	2.00	0.44
33:BL:30:THR:HG22	33:BL:36:LYS:HZ2	1.79	0.44
33:BL:38:GLN:CG	33:BL:39:LYS:H	2.26	0.44
33:BL:56:PRO:HG3	33:BL:60:ARG:HH21	1.82	0.44
35:BN:48:VAL:O	35:BN:52:ILE:HG12	2.18	0.44
23:BB:2335:A:OP2	36:BO:9:ARG:HA	2.18	0.44
37:BP:38:ARG:CG	37:BP:39:LEU:H	2.28	0.44
37:BP:54:LEU:HB3	37:BP:55:HIS:H	1.53	0.44
37:BP:23:ASP:H	37:BP:93:LYS:HZ1	1.66	0.44
39:BR:39:LEU:O	39:BR:54:VAL:HG11	2.17	0.44
39:BR:43:ASN:OD1	39:BR:56:GLY:HA3	2.18	0.44
39:BR:77:PHE:O	39:BR:78:ARG:HG3	2.17	0.44
41:BT:31:VAL:O	41:BT:32:LEU:O	2.36	0.44
42:BU:12:VAL:CG1	42:BU:18:LYS:H	2.31	0.44
23:BB:850:U:O2'	45:BY:21:ALA:HA	2.18	0.44
1:CA:1143:G:O2'	1:CA:1144:G:H5'	2.16	0.44
1:CA:1333:A:H2'	1:CA:1334:G:O4'	2.17	0.44
1:CA:1464:U:H2'	1:CA:1465:A:H8	1.83	0.44
1:CA:1432:G:H1'	1:CA:1468:A:H61	1.81	0.44
1:CA:182:A:O2'	1:CA:183:C:O5'	2.34	0.44
1:CA:28:A:O2'	1:CA:296:U:OP1	2.32	0.44
1:CA:41:G:H2'	1:CA:42:G:H8	1.82	0.44
1:CA:541:G:H2'	1:CA:542:G:C8	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:551:U:H2'	1:CA:552:U:C6	2.53	0.44
1:CA:807:A:N3	25:DC:4:LYS:HE3	2.31	0.44
20:CB:21:TYR:O	20:CB:22:TRP:O	2.36	0.44
3:CD:84:ASN:HD21	3:CD:87:GLU:HG3	1.82	0.44
4:CE:82:HIS:CE1	4:CE:146:MET:HB2	2.53	0.44
6:CG:108:ARG:HH11	6:CG:108:ARG:HG3	1.83	0.44
7:CH:10:LEU:HD22	7:CH:74:ILE:CD1	2.43	0.44
1:CA:878:A:H5''	7:CH:80:PRO:HG2	2.00	0.44
9:CJ:37:ARG:N	9:CJ:76:ILE:HG12	2.33	0.44
11:CL:100:ALA:O	11:CL:101:LEU:C	2.56	0.44
1:CA:523:A:N1	11:CL:88:ASP:HB2	2.33	0.44
19:CT:73:ARG:O	19:CT:74:HIS:C	2.55	0.44
23:DB:1252:G:H5''	55:DB:3462:HOH:O	2.18	0.44
23:DB:1311:G:H21	23:DB:1603:A:H62	1.65	0.44
23:DB:153:U:H2'	23:DB:154:U:C6	2.53	0.44
23:DB:2313:C:H2'	23:DB:2314:A:C8	2.53	0.44
23:DB:2061:G:H5''	23:DB:2503:A:N1	2.32	0.44
23:DB:2556:C:O2'	23:DB:2557:G:H5'	2.17	0.44
23:DB:2849:U:C4'	23:DB:2868:A:C2	3.01	0.44
23:DB:326:G:O2'	23:DB:327:G:H5'	2.17	0.44
23:DB:844:A:H2	23:DB:845:A:N1	2.16	0.44
23:DB:930:G:H5'	23:DB:931:U:P	2.57	0.44
23:DB:998:C:H2'	23:DB:999:U:O4'	2.17	0.44
25:DC:182:LYS:HE2	25:DC:264:LYS:HZ2	1.83	0.44
26:DD:181:ASP:OD1	26:DD:184:ARG:HB3	2.17	0.44
26:DD:180:VAL:HA	26:DD:187:LEU:HA	1.99	0.44
27:DE:139:LYS:HA	27:DE:143:LEU:HD23	1.98	0.44
28:DF:120:SER:HB3	28:DF:127:TYR:HE1	1.83	0.44
28:DF:8:LYS:HA	28:DF:12:VAL:CG2	2.48	0.44
30:DH:45:GLU:O	30:DH:45:GLU:HG3	2.18	0.44
52:DI:4:VAL:HG22	52:DI:5:GLN:N	2.33	0.44
31:DJ:73:VAL:CG2	31:DJ:74:TYR:N	2.81	0.44
33:DL:101:ILE:HG22	33:DL:101:ILE:O	2.17	0.44
33:DL:19:LEU:O	33:DL:21:ARG:N	2.50	0.44
33:DL:59:ARG:HH21	50:D3:5:THR:CB	2.30	0.44
23:DB:630:G:H1	33:DL:69:ARG:HH12	1.64	0.44
34:DM:16:ARG:O	34:DM:17:ASN:HB3	2.17	0.44
34:DM:4:PRO:HB2	34:DM:69:PRO:HG3	1.99	0.44
35:DN:103:ARG:HB3	35:DN:107:ASN:N	2.33	0.44
36:DO:29:HIS:CB	36:DO:36:TYR:HB2	2.48	0.44
38:DQ:111:LYS:HE2	39:DR:52:PRO:CG	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:61:LEU:HB3	41:DT:62:VAL:H	1.55	0.44
41:DT:64:LYS:N	41:DT:64:LYS:HE3	2.33	0.44
41:DT:76:ARG:HG2	41:DT:77:ARG:H	1.83	0.44
45:DY:36:GLU:HG2	45:DY:37:ARG:N	2.32	0.44
1:AA:1129:C:O2'	1:AA:1139:G:N7	2.42	0.44
1:AA:1248:A:H2	8:AI:71:ILE:HD11	1.82	0.44
1:AA:1368:A:O2'	1:AA:1369:C:H5'	2.18	0.44
1:AA:1483:A:H2	23:BB:1959:G:HO2'	1.64	0.44
1:AA:1520:C:H2'	1:AA:1521:C:C6	2.53	0.44
1:AA:542:G:O2'	1:AA:543:U:H5'	2.18	0.44
1:AA:551:U:H2'	1:AA:552:U:C6	2.53	0.44
20:AB:58:LYS:CG	20:AB:62:ARG:HH21	2.26	0.44
3:AD:66:VAL:CG1	3:AD:70:GLN:HB3	2.47	0.44
8:AI:33:SER:C	8:AI:35:GLU:H	2.21	0.44
21:AU:31:VAL:O	21:AU:32:ARG:C	2.56	0.44
48:B1:42:VAL:HG12	48:B1:43:ARG:N	2.33	0.44
50:B3:57:VAL:HG13	50:B3:58:ILE:N	2.32	0.44
22:BA:107:G:O2'	22:BA:108:A:H5'	2.18	0.44
22:BA:54:G:O2'	22:BA:55:U:H5'	2.17	0.44
22:BA:78:A:H61	22:BA:98:G:H1'	1.83	0.44
23:BB:1072:C:N3	23:BB:1092:C:N4	2.65	0.44
23:BB:106:C:H2'	23:BB:107:G:H8	1.83	0.44
23:BB:1157:G:H2'	23:BB:1158:C:C6	2.53	0.44
23:BB:1446:C:H2'	23:BB:1447:C:C6	2.52	0.44
23:BB:2073:C:O2'	23:BB:2074:U:H5'	2.18	0.44
23:BB:2438:U:O2'	23:BB:2439:A:H5''	2.18	0.44
23:BB:2533:U:H2'	23:BB:2534:A:O4'	2.17	0.44
23:BB:465:G:H2'	23:BB:466:A:C8	2.52	0.44
23:BB:499:U:H2'	23:BB:500:G:O4'	2.17	0.44
23:BB:500:G:H1'	23:BB:505:A:N6	2.33	0.44
23:BB:539:G:H2'	23:BB:540:C:C6	2.53	0.44
25:BC:225:ASN:O	25:BC:227:VAL:N	2.51	0.44
27:BE:97:ASN:HB2	27:BE:100:MET:HB2	1.99	0.44
30:BH:62:LEU:O	30:BH:66:ASN:N	2.51	0.44
52:BI:19:PRO:HG2	52:BI:22:PRO:HB2	2.00	0.44
52:BI:72:THR:HB	52:BI:73:PRO:HD2	1.99	0.44
31:BJ:37:ARG:HH21	31:BJ:110:PRO:HG3	1.83	0.44
33:BL:122:VAL:HG22	33:BL:139:GLY:O	2.18	0.44
33:BL:88:GLY:C	33:BL:90:VAL:H	2.20	0.44
23:BB:872:U:H1'	34:BM:68:PHE:CZ	2.53	0.44
37:BP:48:ALA:C	37:BP:49:ILE:HG13	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BP:83:ILE:HG22	37:BP:84:SER:N	2.33	0.44
39:BR:85:LYS:C	39:BR:87:GLN:N	2.71	0.44
23:BB:992:C:H1'	39:BR:91:GLN:HB2	1.99	0.44
40:BS:31:GLN:C	40:BS:33:LEU:N	2.71	0.44
40:BS:48:LYS:O	40:BS:52:GLU:HG3	2.18	0.44
41:BT:62:VAL:O	41:BT:63:VAL:O	2.35	0.44
41:BT:68:LYS:HE3	41:BT:73:ARG:HG3	1.99	0.44
43:BW:34:SER:OG	43:BW:39:GLN:NE2	2.51	0.44
1:CA:372:C:C1'	1:CA:373:A:OP2	2.66	0.44
1:CA:57:G:N2	1:CA:388:G:C6	2.86	0.44
1:CA:56:U:O5'	1:CA:56:U:H6	2.00	0.44
1:CA:649:A:H2'	1:CA:650:G:O4'	2.17	0.44
20:CB:42:LEU:C	20:CB:44:LYS:H	2.20	0.44
3:CD:197:HIS:O	3:CD:200:VAL:HG12	2.17	0.44
1:CA:16:A:O2'	4:CE:20:VAL:HG23	2.18	0.44
5:CF:99:ALA:O	5:CF:100:SER:HB2	2.18	0.44
8:CI:62:LEU:HD22	8:CI:62:LEU:H	1.83	0.44
10:CK:82:GLU:CB	10:CK:108:ASN:HB3	2.48	0.44
11:CL:20:VAL:HA	11:CL:94:TYR:HE2	1.83	0.44
13:CN:55:SER:OG	13:CN:56:PRO:HD2	2.18	0.44
1:CA:617:G:C5'	15:CP:46:LYS:HD3	2.48	0.44
16:CQ:23:ALA:HB1	16:CQ:40:THR:CG2	2.48	0.44
16:CQ:68:LYS:C	16:CQ:70:LYS:H	2.21	0.44
48:D1:49:LYS:HZ2	48:D1:50:GLU:N	2.15	0.44
22:DA:22:U:H2'	22:DA:23:G:H8	1.75	0.44
22:DA:29:A:H8	22:DA:29:A:OP1	2.01	0.44
23:DB:1057:A:C8	23:DB:1086:A:C8	3.05	0.44
23:DB:1157:G:N3	45:DY:10:ARG:NH2	2.66	0.44
23:DB:138:U:O3'	23:DB:139:U:H3'	2.17	0.44
23:DB:1547:C:H2'	23:DB:1548:A:H8	1.81	0.44
23:DB:1710:G:H2'	23:DB:1711:A:H8	1.83	0.44
23:DB:1965:C:H3'	23:DB:1966:A:H8	1.83	0.44
23:DB:2201:G:O2'	23:DB:2202:U:H5'	2.17	0.44
23:DB:2604:U:O2'	23:DB:2605:U:H5'	2.17	0.44
23:DB:2643:G:H2'	23:DB:2644:G:O4'	2.18	0.44
23:DB:2787:C:H2'	23:DB:2788:C:C6	2.53	0.44
23:DB:2811:G:H2'	23:DB:2812:G:H8	1.83	0.44
23:DB:502:A:N6	23:DB:505:A:N1	2.66	0.44
25:DC:22:GLU:HB2	25:DC:202:ARG:HG3	1.99	0.44
26:DD:5:VAL:HG21	26:DD:28:GLU:HA	2.00	0.44
26:DD:69:ALA:HB2	26:DD:91:THR:OG1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:89:GLU:HG2	26:DD:93:GLY:CA	2.41	0.44
27:DE:126:VAL:HG13	27:DE:126:VAL:O	2.18	0.44
27:DE:151:GLY:O	27:DE:158:PHE:HE1	2.01	0.44
23:DB:323:C:C5	27:DE:165:HIS:HB2	2.53	0.44
27:DE:168:ASP:OD2	27:DE:170:ARG:HG2	2.18	0.44
28:DF:107:VAL:CG1	28:DF:175:PRO:HG3	2.48	0.44
29:DG:53:PRO:HG2	29:DG:61:TRP:CE3	2.52	0.44
29:DG:72:ASN:O	29:DG:76:ILE:HG13	2.18	0.44
30:DH:134:VAL:HG12	30:DH:134:VAL:O	2.18	0.44
30:DH:27:ARG:HH11	46:DZ:56:ARG:HD2	1.82	0.44
30:DH:31:VAL:CB	30:DH:32:PRO:HD3	2.38	0.44
52:DI:21:PRO:HB2	52:DI:22:PRO:CD	2.43	0.44
34:DM:133:LYS:CG	34:DM:134:THR:H	2.31	0.44
35:DN:112:TYR:HE2	47:D0:43:THR:HB	1.82	0.44
35:DN:32:GLU:HG3	35:DN:33:ILE:H	1.80	0.44
40:DS:17:VAL:HG11	40:DS:103:ILE:CD1	2.48	0.44
40:DS:4:ILE:HA	40:DS:106:VAL:HG12	1.98	0.44
41:DT:31:VAL:HA	41:DT:84:TYR:HA	2.00	0.44
41:DT:76:ARG:N	41:DT:76:ARG:HD3	2.33	0.44
41:DT:55:VAL:HG21	41:DT:86:THR:H	1.82	0.44
42:DU:4:ILE:HD13	42:DU:4:ILE:C	2.38	0.44
42:DU:9:GLU:HB3	42:DU:21:ARG:HG2	1.99	0.44
44:DX:1:MET:N	44:DX:6:LEU:HG	2.33	0.44
1:AA:1378:C:H2'	1:AA:1378:C:O2	2.16	0.43
1:AA:234:C:H2'	1:AA:235:C:C6	2.53	0.43
1:AA:847:G:H2'	1:AA:848:C:C6	2.52	0.43
1:AA:960:U:H1'	1:AA:961:U:OP2	2.17	0.43
20:AB:112:ARG:HG3	20:AB:112:ARG:NH1	2.32	0.43
3:AD:187:ARG:HG3	3:AD:188:SER:N	2.32	0.43
4:AE:114:LEU:C	4:AE:116:VAL:H	2.21	0.43
5:AF:51:ILE:CD1	5:AF:86:ARG:HG3	2.48	0.43
6:AG:114:SER:O	6:AG:118:ARG:HG3	2.18	0.43
6:AG:142:ARG:HB2	6:AG:142:ARG:CZ	2.48	0.43
7:AH:92:PRO:HA	7:AH:93:LYS:NZ	2.32	0.43
8:AI:97:LEU:N	8:AI:97:LEU:HD22	2.33	0.43
10:AK:35:ASP:CB	10:AK:41:LEU:HD11	2.48	0.43
13:AN:20:PHE:HA	13:AN:24:ALA:CB	2.46	0.43
11:AL:6:LEU:HB3	16:AQ:33:TYR:CZ	2.52	0.43
18:AS:51:HIS:C	18:AS:53:GLY:H	2.21	0.43
23:BB:1599:U:C2	23:BB:1600:C:C5	3.06	0.43
23:BB:1658:C:OP1	26:BD:136:ASN:CA	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1659:G:C6	23:BB:2002:G:C6	3.06	0.43
23:BB:1775:U:H2'	23:BB:1776:G:C5'	2.48	0.43
23:BB:1779:U:OP2	23:BB:1784:A:N6	2.46	0.43
23:BB:1914:C:H2'	23:BB:1915:U:O4'	2.17	0.43
23:BB:2061:G:H3'	55:BB:3546:HOH:O	2.18	0.43
23:BB:2385:C:H2'	23:BB:2386:A:C8	2.53	0.43
23:BB:2886:A:C2	23:BB:2887:A:N7	2.85	0.43
23:BB:30:G:H2'	23:BB:31:C:C6	2.53	0.43
23:BB:855:G:H21	43:BW:23:LYS:CG	2.30	0.43
23:BB:728:G:H4'	25:BC:12:ARG:NH2	2.33	0.43
25:BC:128:THR:CA	25:BC:190:THR:HA	2.41	0.43
25:BC:19:VAL:CG1	25:BC:20:ASN:H	2.20	0.43
25:BC:260:LYS:C	25:BC:262:THR:H	2.21	0.43
23:BB:2052:A:C5'	26:BD:148:GLN:O	2.65	0.43
26:BD:32:ASN:OD1	26:BD:73:VAL:HB	2.18	0.43
28:BF:136:ILE:C	28:BF:138:PRO:HD3	2.39	0.43
29:BG:102:ILE:CG2	29:BG:104:LEU:HD11	2.48	0.43
29:BG:97:VAL:HG23	29:BG:124:CYS:SG	2.58	0.43
52:BI:83:ALA:CA	52:BI:100:ILE:HD11	2.48	0.43
33:BL:115:GLU:OE1	33:BL:115:GLU:HA	2.17	0.43
23:BB:661:A:O2'	33:BL:23:ILE:HG12	2.18	0.43
34:BM:80:VAL:HG12	34:BM:81:ARG:HG2	2.00	0.43
34:BM:8:LYS:O	34:BM:9:PHE:CD1	2.71	0.43
37:BP:38:ARG:O	37:BP:39:LEU:HB2	2.18	0.43
39:BR:46:GLU:O	39:BR:47:VAL:CB	2.66	0.43
39:BR:37:GLU:OE1	39:BR:62:GLU:HB2	2.18	0.43
39:BR:90:ARG:NH1	39:BR:91:GLN:N	2.64	0.43
39:BR:97:LYS:HA	39:BR:97:LYS:HD2	1.82	0.43
42:BU:90:LYS:CD	42:BU:91:LYS:H	2.32	0.43
24:BV:4:ILE:HD11	24:BV:61:LEU:HD12	2.00	0.43
24:BV:79:ARG:H	24:BV:79:ARG:HG2	1.46	0.43
44:BX:10:SER:C	44:BX:12:GLU:H	2.22	0.43
44:BX:22:LEU:CA	44:BX:47:ARG:HH22	2.31	0.43
23:BB:2231:U:OP1	46:BZ:28:VAL:HG12	2.17	0.43
46:BZ:59:ARG:HG2	46:BZ:62:LYS:HB2	1.99	0.43
1:CA:1125:U:HO2'	1:CA:1126:U:P	2.41	0.43
1:CA:1211:U:H1'	1:CA:1213:A:C2	2.53	0.43
1:CA:1463:U:H2'	1:CA:1464:U:C6	2.53	0.43
1:CA:236:A:H2'	1:CA:237:G:C8	2.53	0.43
1:CA:255:G:H2'	1:CA:256:U:C6	2.53	0.43
20:CB:38:HIS:O	20:CB:39:ILE:HD13	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:168:ARG:HE	2:CC:169:GLU:CA	2.30	0.43
3:CD:121:ALA:CA	3:CD:145:ARG:HG3	2.48	0.43
4:CE:88:HIS:CE1	4:CE:137:ARG:HH11	2.35	0.43
5:CF:48:ALA:H	17:CR:65:SER:HB2	1.82	0.43
8:CI:27:ILE:CG2	8:CI:34:LEU:HB2	2.31	0.43
9:CJ:77:VAL:O	9:CJ:78:GLU:C	2.56	0.43
10:CK:113:THR:CG2	21:CU:28:LEU:HD21	2.48	0.43
11:CL:93:ARG:N	11:CL:93:ARG:HD2	2.29	0.43
14:CO:2:LEU:O	14:CO:3:SER:HB3	2.18	0.43
18:CS:30:LEU:HB2	18:CS:48:ILE:HA	2.00	0.43
18:CS:39:ILE:HB	18:CS:66:VAL:O	2.18	0.43
48:D1:27:ARG:HE	48:D1:27:ARG:H	1.65	0.43
23:DB:1000:A:C8	23:DB:1154:G:N2	2.86	0.43
23:DB:1346:G:C6	23:DB:1601:G:C6	3.05	0.43
23:DB:1370:C:H2'	23:DB:1371:G:O4'	2.18	0.43
23:DB:1418:G:N2	23:DB:1579:A:C8	2.86	0.43
23:DB:1634:A:H3'	23:DB:1635:A:C5'	2.48	0.43
23:DB:2662:A:H2'	23:DB:2663:G:O4'	2.18	0.43
23:DB:722:A:H2'	23:DB:723:C:O4'	2.18	0.43
23:DB:946:C:H2'	23:DB:947:A:C8	2.48	0.43
25:DC:179:GLU:CB	25:DC:266:ILE:HA	2.48	0.43
23:DB:2787:C:H5'	26:DD:65:ALA:HB3	2.00	0.43
26:DD:33:ARG:HD2	26:DD:89:GLU:OE2	2.17	0.43
27:DE:129:PRO:O	27:DE:130:LYS:HB2	2.18	0.43
27:DE:147:LEU:HB3	27:DE:167:VAL:CG1	2.45	0.43
27:DE:192:ALA:HB1	27:DE:199:MET:HG2	2.00	0.43
27:DE:73:ILE:O	27:DE:73:ILE:HG22	2.18	0.43
28:DF:141:ASP:C	28:DF:143:ASP:H	2.21	0.43
31:DJ:41:LYS:HG2	38:DQ:63:ARG:CZ	2.48	0.43
23:DB:2250:G:N7	34:DM:82:MET:SD	2.91	0.43
34:DM:97:GLN:H	34:DM:98:PRO:HD3	1.83	0.43
23:DB:2293:G:C5'	36:DO:9:ARG:HH21	2.31	0.43
37:DP:100:ARG:N	37:DP:100:ARG:HD2	2.33	0.43
37:DP:49:ILE:HA	37:DP:62:LYS:O	2.18	0.43
40:DS:77:ASP:O	40:DS:101:SER:HB2	2.18	0.43
24:DV:92:VAL:O	24:DV:92:VAL:HG13	2.18	0.43
1:AA:1099:G:H2'	1:AA:1100:C:O4'	2.18	0.43
1:AA:1424:U:H2'	1:AA:1425:U:H6	1.82	0.43
1:AA:1437:A:O2'	1:AA:1438:G:H5'	2.18	0.43
1:AA:236:A:H2'	1:AA:237:G:C8	2.53	0.43
1:AA:453:G:C6	1:AA:454:G:C6	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AD:169:TRP:HB3	3:AD:183:ARG:HH21	1.84	0.43
3:AD:6:PRO:HG3	3:AD:9:LYS:HD3	1.99	0.43
4:AE:105:ILE:O	4:AE:123:LEU:HA	2.18	0.43
6:AG:19:SER:HB2	6:AG:21:LEU:CD2	2.48	0.43
8:AI:32:ARG:HG2	8:AI:32:ARG:HH11	1.84	0.43
8:AI:38:PHE:O	8:AI:39:GLY:C	2.55	0.43
9:AJ:51:VAL:O	9:AJ:63:ASP:N	2.51	0.43
10:AK:15:VAL:HG21	10:AK:41:LEU:CD1	2.49	0.43
11:AL:26:CYS:HB2	11:AL:29:LYS:HD3	2.01	0.43
13:AN:27:LYS:C	13:AN:27:LYS:HD2	2.38	0.43
19:AT:50:PHE:HE2	19:AT:75:LYS:HA	1.83	0.43
21:AU:37:TYR:C	21:AU:40:PRO:HD2	2.38	0.43
10:AK:110:THR:CG2	21:AU:4:LYS:HA	2.47	0.43
22:BA:32:U:H4'	22:BA:52:A:N6	2.33	0.43
23:BB:1317:G:H2'	23:BB:1318:U:O4'	2.18	0.43
23:BB:1341:G:C2'	23:BB:1397:U:O2'	2.66	0.43
23:BB:1563:U:O2'	23:BB:1564:C:H5'	2.18	0.43
23:BB:1683:U:O2'	23:BB:1684:G:H5'	2.18	0.43
23:BB:1829:A:H3'	23:BB:1830:C:H6	1.83	0.43
23:BB:2293:G:H2'	23:BB:2294:G:O4'	2.18	0.43
23:BB:396:G:O2'	46:BZ:28:VAL:HG11	2.18	0.43
23:BB:409:G:O2'	23:BB:410:G:H5'	2.18	0.43
23:BB:562:U:C4	23:BB:2036:C:O4'	2.71	0.43
23:BB:850:U:H2'	23:BB:851:C:H6	1.82	0.43
23:BB:877:A:H3'	23:BB:899:A:H61	1.82	0.43
25:BC:68:ARG:NH1	25:BC:127:ASN:OD1	2.51	0.43
25:BC:134:ILE:HA	25:BC:135:PRO:HD3	1.91	0.43
25:BC:12:ARG:CZ	25:BC:18:VAL:HG11	2.48	0.43
23:BB:1820:U:O2	25:BC:200:MET:HB2	2.17	0.43
25:BC:222:THR:HA	25:BC:231:HIS:CE1	2.53	0.43
26:BD:132:ALA:O	26:BD:136:ASN:N	2.51	0.43
26:BD:122:VAL:HB	26:BD:141:ARG:HH12	1.82	0.43
26:BD:3:GLY:CA	26:BD:49:GLN:HE22	2.32	0.43
27:BE:3:LEU:HD21	27:BE:109:LEU:HD11	2.00	0.43
27:BE:40:ARG:HD2	27:BE:43:THR:OG1	2.19	0.43
27:BE:75:SER:C	27:BE:77:ILE:H	2.22	0.43
28:BF:14:LYS:HG3	28:BF:15:LEU:N	2.32	0.43
28:BF:169:LEU:O	28:BF:173:ASP:OD1	2.36	0.43
28:BF:67:THR:O	28:BF:83:PRO:HA	2.18	0.43
29:BG:148:ARG:CZ	29:BG:152:ARG:HB2	2.48	0.43
30:BH:142:VAL:HG12	30:BH:143:ILE:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:BH:72:ILE:HD11	30:BH:75:LEU:HD12	1.99	0.43
52:BI:32:VAL:HG13	52:BI:66:PHE:CD2	2.53	0.43
32:BK:12:ASP:OD2	32:BK:85:VAL:HA	2.19	0.43
33:BL:40:SER:O	33:BL:41:ARG:HB2	2.17	0.43
35:BN:35:LYS:CA	35:BN:112:TYR:HB3	2.48	0.43
26:BD:114:LYS:HD3	35:BN:3:HIS:CE1	2.54	0.43
37:BP:104:GLY:O	37:BP:105:LYS:HB3	2.18	0.43
40:BS:33:LEU:HD22	40:BS:33:LEU:N	2.33	0.43
41:BT:36:LYS:O	41:BT:81:LYS:HD2	2.17	0.43
1:CA:1051:C:H2'	1:CA:1052:U:C6	2.54	0.43
1:CA:1092:A:H5''	6:CG:3:ARG:NH1	2.33	0.43
1:CA:1236:A:H4'	1:CA:1304:G:H4'	2.01	0.43
1:CA:488:C:H2'	1:CA:489:C:C6	2.52	0.43
1:CA:695:A:H2'	1:CA:696:A:C8	2.53	0.43
1:CA:908:A:O2'	1:CA:909:A:H5'	2.18	0.43
1:CA:1080:A:C4'	4:CE:20:VAL:HG21	2.43	0.43
5:CF:19:PRO:HG2	5:CF:20:GLY:H	1.83	0.43
8:CI:79:ARG:HA	8:CI:82:ILE:HD12	1.99	0.43
16:CQ:64:ARG:HG3	16:CQ:65:PRO:HD2	1.98	0.43
10:CK:121:ARG:HH21	21:CU:34:ARG:CD	2.30	0.43
47:D0:29:VAL:HG22	47:D0:30:ASP:H	1.81	0.43
51:D4:2:LYS:HZ3	51:D4:2:LYS:HB2	1.82	0.43
23:DB:1221:C:H2'	23:DB:1222:U:C6	2.53	0.43
23:DB:515:A:H2	23:DB:1260:A:N3	2.16	0.43
23:DB:1261:C:C2'	23:DB:1262:A:O5'	2.67	0.43
23:DB:1509:A:C4'	23:DB:1510:G:H5'	2.48	0.43
23:DB:1561:C:H2'	23:DB:1562:U:H6	1.81	0.43
23:DB:2457:U:C2'	23:DB:2458:G:H5'	2.48	0.43
23:DB:268:C:O2	23:DB:268:C:H2'	2.17	0.43
23:DB:285:G:C5	23:DB:356:G:C6	3.06	0.43
23:DB:319:G:O2'	23:DB:320:A:H5'	2.18	0.43
23:DB:557:C:O3'	31:DJ:113:PRO:HB2	2.18	0.43
25:DC:42:ARG:CA	25:DC:42:ARG:NE	2.81	0.43
27:DE:149:ILE:HD11	27:DE:188:MET:H	1.81	0.43
27:DE:141:MET:HG3	27:DE:185:LYS:HE3	1.99	0.43
27:DE:49:ARG:HB2	27:DE:51:GLU:CD	2.38	0.43
28:DF:140:ILE:C	28:DF:142:TYR:H	2.21	0.43
52:DI:23:VAL:HG12	52:DI:24:GLY:H	1.83	0.43
52:DI:73:PRO:HA	52:DI:74:PRO:HD3	1.92	0.43
31:DJ:24:THR:O	31:DJ:28:LEU:HG	2.19	0.43
34:DM:9:PHE:CE1	34:DM:11:LYS:HB2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DN:126:ALA:C	35:DN:127:GLU:HG2	2.39	0.43
23:DB:1010:A:H5'	38:DQ:61:ILE:CG2	2.48	0.43
39:DR:58:VAL:O	39:DR:59:ILE:HB	2.17	0.43
39:DR:92:TRP:CE3	39:DR:93:PHE:N	2.87	0.43
1:AA:1113:C:O2'	1:AA:1114:C:H5'	2.17	0.43
1:AA:113:G:N2	1:AA:353:A:H8	2.16	0.43
1:AA:12:U:H4'	1:AA:526:C:H4'	1.99	0.43
1:AA:177:G:N3	1:AA:177:G:O4'	2.51	0.43
1:AA:179:A:H2'	1:AA:180:U:C6	2.54	0.43
1:AA:237:G:O2'	1:AA:238:A:H5'	2.18	0.43
1:AA:429:U:H4'	1:AA:430:A:O5'	2.16	0.43
1:AA:57:G:N2	1:AA:388:G:C6	2.86	0.43
1:AA:605:U:H2'	1:AA:606:G:C8	2.53	0.43
1:AA:729:A:H2'	1:AA:730:G:H8	1.83	0.43
1:AA:90:C:H2'	1:AA:91:U:H6	1.83	0.43
2:AC:166:TRP:HB3	2:AC:167:TYR:H	1.45	0.43
2:AC:21:TRP:CD1	2:AC:58:ARG:HD2	2.53	0.43
3:AD:194:ILE:HD11	3:AD:199:ILE:HD11	2.00	0.43
3:AD:51:GLY:O	3:AD:55:ARG:HB2	2.18	0.43
4:AE:95:MET:CE	4:AE:114:LEU:HD21	2.48	0.43
11:AL:38:THR:CG2	11:AL:48:LEU:HD12	2.48	0.43
14:AO:38:LEU:HB3	14:AO:42:PHE:HE1	1.83	0.43
16:AQ:32:ILE:HG13	16:AQ:33:TYR:N	2.33	0.43
16:AQ:60:ILE:HD13	16:AQ:60:ILE:H	1.83	0.43
18:AS:41:PRO:HG2	18:AS:42:ASN:H	1.83	0.43
18:AS:7:GLY:O	18:AS:9:PHE:N	2.51	0.43
47:B0:51:ARG:NH1	47:B0:56:LYS:O	2.51	0.43
51:B4:31:PRO:O	51:B4:32:LYS:C	2.56	0.43
23:BB:1290:C:O2'	23:BB:1291:C:H5'	2.19	0.43
23:BB:135:U:C2'	23:BB:136:G:C8	2.92	0.43
23:BB:1680:U:H2'	23:BB:1681:G:O4'	2.18	0.43
23:BB:1710:G:H2'	23:BB:1711:A:H8	1.84	0.43
23:BB:1812:U:H2'	23:BB:1813:G:C8	2.53	0.43
23:BB:2014:A:OP1	40:BS:95:ARG:NE	2.39	0.43
23:BB:2278:A:H62	43:BW:10:ARG:HG2	1.82	0.43
23:BB:2289:G:O2'	23:BB:2290:G:H5'	2.18	0.43
23:BB:2355:G:H4'	43:BW:20:LEU:CD2	2.45	0.43
23:BB:2479:U:H6	23:BB:2479:U:O5'	2.01	0.43
23:BB:255:A:H2'	23:BB:256:A:O4'	2.18	0.43
23:BB:2564:A:OP1	23:BB:2648:G:H4'	2.18	0.43
23:BB:2728:U:H2'	23:BB:2729:G:H8	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2866:U:O3'	23:BB:2867:G:H4'	2.18	0.43
23:BB:540:C:H2'	23:BB:541:A:H8	1.83	0.43
23:BB:711:G:H2'	23:BB:712:G:O4'	2.17	0.43
23:BB:719:C:O2'	23:BB:720:U:H5'	2.18	0.43
23:BB:725:G:H2'	23:BB:726:G:C1'	2.48	0.43
23:BB:975:A:N6	23:BB:989:G:O2'	2.52	0.43
23:BB:784:G:H5''	25:BC:225:ASN:ND2	2.33	0.43
25:BC:33:LEU:HB3	25:BC:34:GLU:H	1.72	0.43
25:BC:74:PRO:HB2	25:BC:96:LYS:CG	2.36	0.43
25:BC:78:GLU:HB2	25:BC:92:LEU:HB3	1.99	0.43
26:BD:103:ASP:HA	26:BD:105:LYS:HE2	1.99	0.43
27:BE:27:LEU:CD2	27:BE:100:MET:HG2	2.48	0.43
27:BE:127:GLU:OE1	27:BE:130:LYS:HG3	2.19	0.43
27:BE:139:LYS:HE2	27:BE:139:LYS:CA	2.48	0.43
29:BG:169:ARG:C	29:BG:171:LYS:N	2.72	0.43
30:BH:124:THR:HG22	30:BH:125:THR:H	1.82	0.43
30:BH:114:GLU:HB3	30:BH:134:VAL:HA	1.99	0.43
30:BH:60:GLU:OE1	30:BH:61:VAL:HG23	2.18	0.43
30:BH:66:ASN:C	30:BH:68:ARG:H	2.22	0.43
52:BI:12:VAL:HG23	52:BI:41:PHE:CE2	2.53	0.43
31:BJ:41:LYS:CD	31:BJ:44:TYR:HB2	2.45	0.43
31:BJ:90:GLU:OE1	31:BJ:94:ALA:HB2	2.17	0.43
32:BK:14:SER:O	32:BK:52:VAL:HB	2.18	0.43
32:BK:80:ASP:CG	37:BP:71:ARG:HG2	2.39	0.43
34:BM:67:VAL:HG13	34:BM:67:VAL:O	2.18	0.43
34:BM:98:PRO:HG2	34:BM:99:GLY:H	1.83	0.43
37:BP:22:GLY:HA3	37:BP:91:VAL:HG13	2.00	0.43
37:BP:84:SER:O	37:BP:86:LYS:N	2.50	0.43
38:BQ:26:ALA:C	38:BQ:28:SER:H	2.22	0.43
39:BR:17:GLY:C	39:BR:18:GLN:HG2	2.39	0.43
23:BB:1163:G:N2	39:BR:91:GLN:HG3	2.31	0.43
24:BV:82:TYR:HD1	34:BM:36:VAL:HB	1.83	0.43
43:BW:47:GLY:N	43:BW:67:LYS:HZ3	2.15	0.43
46:BZ:64:PHE:C	46:BZ:67:PRO:HD2	2.39	0.43
1:CA:1187:G:OP1	8:CI:114:LYS:HE2	2.17	0.43
1:CA:1301:U:HO2'	1:CA:1302:C:P	2.36	0.43
1:CA:1458:G:O2'	1:CA:1459:G:H5'	2.19	0.43
1:CA:1460:C:O2'	1:CA:1461:G:H5'	2.19	0.43
1:CA:374:A:H5''	1:CA:452:A:N1	2.33	0.43
1:CA:402:G:H2'	1:CA:403:C:C6	2.53	0.43
1:CA:797:C:OP1	10:CK:125:LYS:HE2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:847:G:H2'	1:CA:848:C:C6	2.53	0.43
2:CC:8:GLY:HA2	2:CC:11:LEU:HG	1.99	0.43
3:CD:179:GLY:O	3:CD:180:THR:HB	2.18	0.43
4:CE:155:LYS:H	4:CE:155:LYS:HG3	1.53	0.43
5:CF:43:GLY:HA2	5:CF:58:HIS:NE2	2.33	0.43
8:CI:97:LEU:O	8:CI:102:PHE:HB2	2.17	0.43
12:CM:56:ARG:HG3	12:CM:56:ARG:NH1	2.30	0.43
14:CO:74:VAL:HG11	16:CQ:83:LEU:OXT	2.18	0.43
50:D3:49:VAL:C	50:D3:51:LYS:H	2.22	0.43
22:DA:114:C:H2'	22:DA:115:A:C8	2.53	0.43
23:DB:1180:U:O2'	23:DB:1181:U:H5'	2.18	0.43
23:DB:1236:G:H2'	23:DB:1237:A:OP2	2.18	0.43
23:DB:1535:A:H5'	23:DB:1536:C:C5	2.52	0.43
23:DB:1736:U:H2'	23:DB:1737:G:C8	2.52	0.43
23:DB:1877:A:O2'	23:DB:1878:G:H5'	2.18	0.43
23:DB:2259:U:N3	23:DB:2260:C:C5	2.86	0.43
23:DB:2281:A:N6	43:DW:3:LYS:HE3	2.32	0.43
23:DB:2310:C:H2'	23:DB:2311:A:H5'	1.99	0.43
23:DB:2526:G:C2	51:D4:2:LYS:HE3	2.53	0.43
23:DB:2733:A:C8	23:DB:2733:A:C3'	3.02	0.43
23:DB:192:C:O2'	23:DB:802:A:N3	2.51	0.43
23:DB:908:C:O2'	23:DB:909:A:H5'	2.18	0.43
23:DB:935:C:O2'	23:DB:936:A:H5'	2.18	0.43
23:DB:947:A:H2'	23:DB:948:C:H6	1.76	0.43
25:DC:21:PRO:CD	25:DC:202:ARG:NH1	2.81	0.43
25:DC:51:ARG:O	25:DC:52:HIS:C	2.57	0.43
26:DD:181:ASP:CG	26:DD:184:ARG:HB3	2.39	0.43
26:DD:18:ASP:C	26:DD:20:VAL:N	2.70	0.43
26:DD:82:PHE:HB3	26:DD:83:ARG:H	1.42	0.43
27:DE:130:LYS:HZ2	27:DE:130:LYS:HB2	1.80	0.43
28:DF:7:TYR:OH	28:DF:29:ARG:HG2	2.18	0.43
23:DB:2530:A:H5'	29:DG:176:LYS:O	2.18	0.43
31:DJ:120:ARG:HB3	31:DJ:121:LYS:CE	2.48	0.43
31:DJ:95:ARG:NE	31:DJ:95:ARG:N	2.65	0.43
32:DK:70:ARG:HG3	32:DK:70:ARG:O	2.18	0.43
33:DL:61:LEU:HB3	33:DL:62:PRO:HD3	1.99	0.43
22:DA:7:G:H4'	36:DO:29:HIS:CD2	2.53	0.43
37:DP:13:LYS:NZ	37:DP:15:ASP:HB2	2.33	0.43
38:DQ:73:ILE:HG13	38:DQ:74:SER:N	2.22	0.43
39:DR:20:VAL:HG12	39:DR:97:LYS:HA	1.99	0.43
39:DR:38:VAL:HA	39:DR:61:ALA:HB3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DS:29:VAL:HG21	40:DS:69:LEU:C	2.38	0.43
44:DX:19:LEU:C	44:DX:21:LEU:H	2.22	0.43
1:AA:1288:A:O2'	1:AA:1289:A:H5'	2.18	0.43
1:AA:195:A:H1'	1:AA:222:C:O2'	2.18	0.43
1:AA:340:U:H2'	1:AA:341:C:C6	2.54	0.43
1:AA:392:C:H2'	1:AA:393:A:C8	2.54	0.43
1:AA:462:G:H3'	1:AA:463:U:C5'	2.48	0.43
20:AB:61:SER:HB3	20:AB:223:GLY:O	2.17	0.43
2:AC:113:LYS:HA	2:AC:113:LYS:HE2	2.00	0.43
2:AC:147:GLY:HA3	2:AC:171:ARG:H	1.84	0.43
1:AA:932:C:H5'	6:AG:3:ARG:NH2	2.34	0.43
7:AH:10:LEU:HA	7:AH:74:ILE:HD11	2.00	0.43
8:AI:10:ARG:HA	8:AI:77:ALA:CB	2.49	0.43
8:AI:85:ALA:O	8:AI:88:GLU:HB2	2.18	0.43
9:AJ:41:PRO:O	9:AJ:71:LEU:HD13	2.18	0.43
9:AJ:39:PRO:CB	9:AJ:74:VAL:HG22	2.48	0.43
11:AL:121:PRO:C	11:AL:123:ALA:H	2.21	0.43
11:AL:41:PRO:CB	11:AL:88:ASP:HB3	2.48	0.43
14:AO:46:LYS:HE3	14:AO:46:LYS:HB2	1.84	0.43
1:AA:130:A:C8	16:AQ:64:ARG:HG3	2.53	0.43
1:AA:323:U:H4'	19:AT:16:ALA:HB3	2.01	0.43
48:B1:36:LYS:HB3	48:B1:37:LYS:H	1.67	0.43
22:BA:91:C:H2'	22:BA:92:C:H6	1.83	0.43
23:BB:1025:G:H1'	23:BB:1135:C:O5'	2.18	0.43
23:BB:1081:U:H4'	52:BI:123:ALA:HB1	2.01	0.43
23:BB:1178:C:C2	23:BB:1179:G:N7	2.87	0.43
23:BB:1183:U:C2	23:BB:1184:U:C5	3.06	0.43
23:BB:1463:C:H2'	23:BB:1464:G:H8	1.82	0.43
23:BB:1930:G:C2'	23:BB:1931:U:OP2	2.67	0.43
23:BB:211:C:O2'	23:BB:212:G:H5'	2.18	0.43
23:BB:2336:A:HO2'	23:BB:2337:G:P	2.41	0.43
23:BB:2457:U:O2'	23:BB:2458:G:H5'	2.18	0.43
23:BB:2733:A:C3'	23:BB:2733:A:C8	3.02	0.43
23:BB:2849:U:C4'	23:BB:2868:A:C2	3.01	0.43
23:BB:507:A:OP1	23:BB:508:A:N7	2.52	0.43
23:BB:709:U:H2'	23:BB:710:U:H6	1.84	0.43
26:BD:140:HIS:HB3	26:BD:141:ARG:H	1.62	0.43
26:BD:164:GLN:HG2	26:BD:165:MET:H	1.83	0.43
27:BE:141:MET:CG	27:BE:143:LEU:HB2	2.48	0.43
27:BE:161:ALA:O	27:BE:162:ARG:HD2	2.18	0.43
27:BE:149:ILE:CD1	27:BE:186:VAL:HG13	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:BF:103:ILE:HG13	28:BF:104:THR:H	1.83	0.43
29:BG:49:LEU:O	29:BG:50:THR:HG23	2.19	0.43
29:BG:57:TYR:CD1	29:BG:57:TYR:N	2.86	0.43
52:BI:77:VAL:HG23	52:BI:78:LEU:N	2.33	0.43
31:BJ:25:LEU:HB2	31:BJ:26:GLY:H	1.57	0.43
31:BJ:64:VAL:HG11	31:BJ:68:LYS:CD	2.34	0.43
33:BL:64:PHE:N	50:B3:24:LYS:HZ2	2.16	0.43
33:BL:82:LEU:HD13	33:BL:83:ALA:H	1.81	0.43
34:BM:64:TRP:CE3	34:BM:102:LEU:HD22	2.54	0.43
35:BN:122:ALA:CB	47:B0:49:ARG:HH12	2.32	0.43
35:BN:24:MET:O	35:BN:28:LEU:HD12	2.18	0.43
35:BN:8:ARG:HB3	35:BN:43:GLU:OE1	2.19	0.43
35:BN:57:THR:HG21	35:BN:61:ALA:HB3	2.01	0.43
36:BO:84:GLU:O	36:BO:87:ILE:HG12	2.18	0.43
37:BP:25:VAL:O	37:BP:25:VAL:HG12	2.18	0.43
37:BP:59:THR:HG1	37:BP:76:HIS:CE1	2.36	0.43
38:BQ:94:LEU:CB	39:BR:13:ARG:HG2	2.49	0.43
39:BR:42:ALA:C	39:BR:44:GLY:N	2.72	0.43
41:BT:68:LYS:HB2	41:BT:68:LYS:HE2	1.82	0.43
45:BY:15:ARG:CA	45:BY:15:ARG:HE	2.31	0.43
1:CA:1248:A:H2'	1:CA:1249:C:H6	1.83	0.43
1:CA:1253:G:N1	1:CA:1285:A:N6	2.66	0.43
1:CA:1521:C:H2'	1:CA:1522:U:C6	2.53	0.43
1:CA:25:C:H2'	1:CA:26:A:C8	2.54	0.43
1:CA:425:G:O2'	1:CA:426:U:H5'	2.19	0.43
1:CA:512:U:O2'	1:CA:513:C:H5'	2.19	0.43
1:CA:521:G:O2'	1:CA:522:C:H5'	2.17	0.43
1:CA:94:G:H4'	1:CA:95:C:H5''	2.01	0.43
20:CB:41:ASN:ND2	20:CB:44:LYS:HB2	2.32	0.43
2:CC:11:LEU:HD22	2:CC:17:TRP:CD1	2.54	0.43
3:CD:33:ILE:O	3:CD:34:GLU:C	2.57	0.43
3:CD:61:ARG:HG3	3:CD:66:VAL:O	2.19	0.43
5:CF:59:TYR:N	5:CF:59:TYR:CD1	2.86	0.43
6:CG:11:ILE:HG22	6:CG:12:LEU:N	2.33	0.43
8:CI:29:ILE:CA	8:CI:64:ILE:HB	2.46	0.43
11:CL:106:VAL:CG1	11:CL:109:ARG:HG3	2.48	0.43
13:CN:80:ARG:HH11	13:CN:80:ARG:HG2	1.84	0.43
14:CO:28:VAL:HB	14:CO:80:LEU:HD11	1.99	0.43
18:CS:19:GLU:O	18:CS:19:GLU:HG3	2.18	0.43
1:CA:1320:C:N4	18:CS:36:ARG:HD2	2.30	0.43
19:CT:15:LYS:HA	19:CT:15:LYS:HD3	1.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:D1:9:LYS:CG	48:D1:24:LYS:HG2	2.49	0.43
48:D1:49:LYS:HD2	48:D1:49:LYS:O	2.17	0.43
23:DB:2361:G:C5'	50:D3:27:ASN:HB2	2.48	0.43
51:D4:10:LEU:HD13	51:D4:25:VAL:HB	2.00	0.43
51:D4:26:ILE:HG23	51:D4:27:CYS:N	2.21	0.43
23:DB:51:G:H1'	23:DB:118:A:H61	1.82	0.43
23:DB:138:U:O3'	23:DB:139:U:H2'	2.19	0.43
23:DB:1826:G:OP1	25:DC:222:THR:HG22	2.18	0.43
23:DB:215:G:H4'	23:DB:216:A:OP1	2.17	0.43
23:DB:2304:G:H22	23:DB:2312:U:H3	1.66	0.43
23:DB:2349:G:OP2	50:D3:41:ARG:HD3	2.17	0.43
23:DB:2660:A:H2'	23:DB:2661:G:O4'	2.17	0.43
23:DB:2727:A:O2'	32:DK:70:ARG:NH2	2.52	0.43
23:DB:743:A:C2'	23:DB:744:U:H5'	2.47	0.43
23:DB:78:U:H2'	23:DB:79:C:H6	1.82	0.43
25:DC:68:ARG:CZ	25:DC:127:ASN:HA	2.46	0.43
26:DD:191:GLY:O	26:DD:192:ALA:HB3	2.17	0.43
26:DD:50:VAL:CG1	26:DD:75:ALA:HB3	2.46	0.43
26:DD:77:ARG:CB	26:DD:77:ARG:HH11	2.30	0.43
34:DM:73:ILE:HG13	34:DM:92:TRP:HB2	2.00	0.43
35:DN:28:LEU:HD23	35:DN:48:VAL:HG11	2.00	0.43
37:DP:86:LYS:C	37:DP:87:ARG:HG3	2.37	0.43
38:DQ:113:LYS:C	38:DQ:115:ALA:H	2.21	0.43
39:DR:1:MET:N	39:DR:47:VAL:HB	2.33	0.43
23:DB:136:G:N1	41:DT:3:ARG:CZ	2.81	0.43
42:DU:81:ARG:HG3	42:DU:92:VAL:HG21	1.99	0.43
23:DB:2281:A:H62	43:DW:3:LYS:HD2	1.83	0.43
45:DY:26:LEU:C	45:DY:28:LEU:N	2.71	0.43
1:AA:109:A:C6	1:AA:326:G:C6	3.06	0.43
1:AA:1213:A:C6	1:AA:1215:G:H1'	2.54	0.43
1:AA:1454:G:H2'	1:AA:1455:G:O4'	2.19	0.43
1:AA:186:C:H2'	1:AA:187:G:O4'	2.18	0.43
1:AA:224:U:H2'	1:AA:225:C:H6	1.82	0.43
1:AA:255:G:H2'	1:AA:256:U:C6	2.53	0.43
1:AA:598:U:H2'	1:AA:599:C:H6	1.82	0.43
1:AA:635:A:H2'	1:AA:636:U:C6	2.53	0.43
20:AB:187:ASP:OD1	20:AB:188:THR:N	2.51	0.43
20:AB:38:HIS:O	20:AB:39:ILE:HD13	2.17	0.43
9:AJ:80:THR:HG22	9:AJ:81:GLU:N	2.29	0.43
15:AP:25:ARG:HD3	15:AP:25:ARG:H	1.83	0.43
18:AS:43:MET:HE1	18:AS:48:ILE:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AU:16:ARG:NH2	21:AU:24:LYS:HE2	2.33	0.43
50:B3:21:PHE:CD1	50:B3:21:PHE:N	2.85	0.43
23:BB:1132:U:C6	31:BJ:85:LYS:HE2	2.53	0.43
23:BB:1201:U:H2'	23:BB:1202:G:H8	1.84	0.43
23:BB:142:A:OP2	23:BB:142:A:H8	2.01	0.43
23:BB:2093:G:H2'	23:BB:2094:A:H8	1.84	0.43
23:BB:2303:G:H2'	23:BB:2304:G:H5'	2.00	0.43
23:BB:234:U:H2'	23:BB:235:U:H6	1.84	0.43
23:BB:2686:G:H2'	23:BB:2687:U:O4'	2.18	0.43
23:BB:440:C:H2'	23:BB:441:U:H6	1.82	0.43
23:BB:466:A:H2'	23:BB:467:G:H5'	2.01	0.43
23:BB:619:G:H2'	23:BB:620:G:H5''	2.00	0.43
23:BB:822:G:H2'	23:BB:823:C:C6	2.53	0.43
23:BB:904:G:H2'	23:BB:905:A:C8	2.52	0.43
23:BB:958:U:H5'	34:BM:18:ARG:HH12	1.83	0.43
25:BC:110:LYS:CE	25:BC:111:ALA:H	2.27	0.43
25:BC:88:ALA:CB	25:BC:156:SER:HB3	2.42	0.43
25:BC:191:LEU:HD22	25:BC:191:LEU:O	2.17	0.43
25:BC:257:ARG:HA	25:BC:257:ARG:CZ	2.47	0.43
25:BC:71:ASP:OD2	25:BC:118:GLY:HA2	2.19	0.43
26:BD:62:LYS:HE3	26:BD:63:PRO:HD2	2.01	0.43
27:BE:109:LEU:HA	27:BE:117:ARG:HB3	2.00	0.43
27:BE:121:VAL:HG12	27:BE:122:GLU:H	1.80	0.43
27:BE:138:LEU:HD22	27:BE:144:GLU:HA	1.99	0.43
27:BE:25:GLU:O	27:BE:26:ALA:C	2.56	0.43
30:BH:62:LEU:HD21	30:BH:137:GLU:HG2	2.01	0.43
30:BH:58:LEU:O	30:BH:62:LEU:HD12	2.19	0.43
52:BI:85:ILE:HD12	52:BI:87:SER:O	2.19	0.43
31:BJ:58:ASN:C	31:BJ:60:ASP:H	2.21	0.43
33:BL:118:THR:N	33:BL:119:PRO:HD3	2.33	0.43
33:BL:39:LYS:HB3	33:BL:46:VAL:HG13	1.99	0.43
34:BM:80:VAL:HG12	34:BM:81:ARG:N	2.33	0.43
37:BP:92:ARG:CZ	37:BP:110:LYS:HE2	2.46	0.43
37:BP:47:ILE:CA	37:BP:63:ILE:HG23	2.48	0.43
39:BR:33:VAL:HB	39:BR:35:PHE:HE2	1.82	0.43
40:BS:43:ALA:O	40:BS:47:VAL:HG13	2.18	0.43
41:BT:15:HIS:HB2	41:BT:31:VAL:HB	1.99	0.43
24:BV:26:PHE:HB2	24:BV:27:PRO:HD2	2.00	0.43
43:BW:15:SER:O	43:BW:16:GLU:HB3	2.17	0.43
46:BZ:49:ARG:O	46:BZ:50:ASP:HB3	2.18	0.43
1:CA:1210:C:H4'	1:CA:1214:C:C4	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1234:C:H2'	1:CA:1235:U:H6	1.83	0.43
1:CA:1280:A:O4'	9:CJ:43:PRO:HG3	2.18	0.43
1:CA:128:G:O2'	1:CA:129:A:H5'	2.18	0.43
1:CA:1348:U:O2	1:CA:1348:U:H2'	2.18	0.43
1:CA:1376:U:H2'	1:CA:1377:A:C8	2.54	0.43
1:CA:206:C:H2'	1:CA:207:C:C6	2.53	0.43
1:CA:239:U:H6	1:CA:239:U:H5''	1.80	0.43
1:CA:113:G:N2	1:CA:353:A:H8	2.16	0.43
1:CA:847:G:H2'	1:CA:848:C:H6	1.82	0.43
1:CA:916:U:H2'	1:CA:917:G:C8	2.52	0.43
20:CB:86:CYS:C	20:CB:88:GLN:H	2.21	0.43
2:CC:137:VAL:HG13	2:CC:148:ILE:HG23	2.00	0.43
3:CD:123:MET:HB2	3:CD:127:ARG:O	2.19	0.43
3:CD:171:GLU:OE2	3:CD:182:LYS:HE3	2.19	0.43
3:CD:64:TYR:CD2	3:CD:93:LEU:HB2	2.54	0.43
4:CE:136:VAL:HG23	4:CE:137:ARG:H	1.83	0.43
8:CI:29:ILE:O	8:CI:32:ARG:HB2	2.18	0.43
17:CR:51:GLN:NE2	17:CR:54:LEU:HB2	2.33	0.43
18:CS:29:PRO:HB3	18:CS:47:THR:HG22	2.01	0.43
19:CT:73:ARG:HG3	19:CT:74:HIS:H	1.83	0.43
21:CU:17:ARG:N	21:CU:17:ARG:HD2	2.12	0.43
48:D1:27:ARG:H	48:D1:27:ARG:NE	2.16	0.43
23:DB:1309:G:C5'	49:D2:8:SER:H	2.30	0.43
50:D3:24:LYS:C	50:D3:26:ALA:H	2.22	0.43
50:D3:25:HIS:CE1	50:D3:46:LYS:HB2	2.53	0.43
51:D4:18:LYS:O	51:D4:19:ARG:HG3	2.17	0.43
22:DA:53:A:C2'	22:DA:54:G:H5'	2.49	0.43
22:DA:13:G:H1'	22:DA:69:G:N2	2.33	0.43
23:DB:1038:G:H2'	23:DB:1039:A:C8	2.52	0.43
23:DB:1616:A:H4'	23:DB:1617:C:OP2	2.19	0.43
23:DB:1774:C:C2'	23:DB:1774:C:O2	2.54	0.43
23:DB:184:C:H2'	23:DB:185:G:C8	2.51	0.43
23:DB:1948:G:C6	23:DB:1959:G:C6	3.06	0.43
23:DB:2072:C:O2'	23:DB:2073:C:H5'	2.19	0.43
23:DB:2096:C:H2'	23:DB:2097:A:C8	2.53	0.43
23:DB:2382:G:O2'	50:D3:38:LYS:HE2	2.18	0.43
23:DB:2617:U:H2'	23:DB:2618:G:C5'	2.47	0.43
23:DB:2691:C:H2'	23:DB:2692:G:H8	1.84	0.43
23:DB:303:G:C6	23:DB:315:G:C6	3.06	0.43
23:DB:79:C:O2'	23:DB:346:A:C8	2.68	0.43
23:DB:511:U:H4'	23:DB:1235:G:H4'	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:585:G:H2'	23:DB:1251:C:N4	2.33	0.43
23:DB:599:A:H2'	23:DB:600:G:H8	1.83	0.43
23:DB:773:U:O2'	25:DC:48:ILE:HG23	2.18	0.43
29:DG:18:ILE:HG13	29:DG:18:ILE:O	2.18	0.43
30:DH:46:PHE:O	30:DH:49:ALA:HB3	2.19	0.43
52:DI:7:TYR:CE1	52:DI:57:VAL:HG11	2.53	0.43
31:DJ:35:ARG:O	31:DJ:52:ASP:HB2	2.18	0.43
31:DJ:95:ARG:HH11	31:DJ:95:ARG:HG3	1.83	0.43
32:DK:19:VAL:HB	32:DK:41:ILE:CG1	2.49	0.43
33:DL:126:ARG:HH11	33:DL:126:ARG:HG2	1.83	0.43
23:DB:588:U:H5'	33:DL:29:LYS:NZ	2.34	0.43
33:DL:54:GLN:CB	33:DL:57:LEU:HD23	2.48	0.43
33:DL:63:LYS:HE3	50:D3:29:ARG:HG3	2.00	0.43
35:DN:103:ARG:NH1	35:DN:106:ASP:OD2	2.51	0.43
35:DN:18:GLN:HB2	35:DN:18:GLN:HE21	1.55	0.43
38:DQ:87:VAL:HB	39:DR:54:VAL:CG2	2.40	0.43
40:DS:58:ALA:HA	40:DS:62:ASP:HB3	2.00	0.43
24:DV:10:LYS:HD2	24:DV:41:GLU:OE2	2.18	0.43
24:DV:57:TYR:CE2	24:DV:77:VAL:HG21	2.53	0.43
24:DV:60:VAL:HG12	24:DV:73:LYS:HG3	2.00	0.43
43:DW:56:HIS:CD2	43:DW:58:LEU:H	2.36	0.43
45:DY:57:GLU:OE1	45:DY:57:GLU:O	2.36	0.43
1:AA:1140:C:O2'	1:AA:1141:C:H5'	2.18	0.43
1:AA:1220:G:H2'	1:AA:1221:G:C8	2.52	0.43
1:AA:1238:A:N3	1:AA:1241:G:H1'	2.34	0.43
1:AA:1343:G:H2'	1:AA:1344:C:C6	2.54	0.43
1:AA:1343:G:H2'	1:AA:1344:C:H6	1.83	0.43
1:AA:1367:C:O2'	1:AA:1368:A:H5'	2.16	0.43
1:AA:65:A:N3	1:AA:65:A:C2'	2.76	0.43
1:AA:695:A:H2'	1:AA:696:A:C8	2.54	0.43
1:AA:77:A:H2'	1:AA:78:A:N7	2.34	0.43
2:AC:139:ASN:HD22	2:AC:139:ASN:N	2.16	0.43
3:AD:64:TYR:N	3:AD:64:TYR:CD1	2.87	0.43
5:AF:32:ALA:CB	5:AF:70:VAL:HG11	2.49	0.43
8:AI:43:ALA:HA	8:AI:44:ARG:NH1	2.34	0.43
8:AI:56:MET:CA	8:AI:59:LYS:HZ2	2.30	0.43
8:AI:66:VAL:CG2	8:AI:74:GLN:HG3	2.48	0.43
12:AM:33:LEU:HB3	12:AM:38:ILE:O	2.18	0.43
21:AU:36:PHE:C	21:AU:40:PRO:HD3	2.38	0.43
23:BB:1115:G:H2'	23:BB:1116:G:H8	1.83	0.43
23:BB:1117:C:C2	23:BB:1118:C:C5	3.07	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1155:A:H61	45:BY:10:ARG:CZ	2.31	0.43
23:BB:161:A:H62	23:BB:165:A:N6	2.10	0.43
23:BB:1877:A:H2'	23:BB:1878:G:C8	2.53	0.43
23:BB:1948:G:C6	23:BB:1959:G:C6	3.06	0.43
23:BB:2184:A:H2'	23:BB:2185:U:C6	2.53	0.43
23:BB:2210:U:N3	23:BB:2212:A:N7	2.65	0.43
23:BB:2728:U:H2'	23:BB:2729:G:C8	2.54	0.43
23:BB:2771:C:O2'	26:BD:173:GLN:NE2	2.52	0.43
23:BB:315:G:H2'	23:BB:316:C:O4'	2.19	0.43
23:BB:636:G:OP2	33:BL:128:THR:HG22	2.18	0.43
25:BC:194:VAL:CG1	25:BC:195:GLY:H	2.17	0.43
25:BC:25:LYS:HA	25:BC:25:LYS:HD2	1.83	0.43
23:BB:2574:G:HO2'	26:BD:151:THR:HG21	1.84	0.43
26:BD:24:VAL:O	26:BD:24:VAL:HG13	2.18	0.43
26:BD:52:THR:HG22	26:BD:76:GLY:H	1.84	0.43
27:BE:150:THR:HB	27:BE:171:ASP:OD1	2.18	0.43
27:BE:149:ILE:O	27:BE:187:VAL:HG22	2.18	0.43
28:BF:3:LEU:HA	28:BF:6:TYR:CG	2.54	0.43
28:BF:45:ASP:OD2	28:BF:83:PRO:HG2	2.18	0.43
30:BH:11:ASN:C	30:BH:13:GLY:H	2.22	0.43
52:BI:15:GLY:O	52:BI:16:MET:HB2	2.16	0.43
31:BJ:35:ARG:NH2	31:BJ:40:HIS:H	2.15	0.43
31:BJ:43:GLU:HG3	31:BJ:44:TYR:CD2	2.53	0.43
23:BB:954:G:OP1	34:BM:17:ASN:HB2	2.19	0.43
34:BM:71:LYS:HD3	34:BM:92:TRP:N	2.33	0.43
37:BP:51:ASN:HA	37:BP:61:ARG:N	2.29	0.43
38:BQ:92:LYS:O	38:BQ:94:LEU:HD22	2.18	0.43
24:BV:42:LEU:HD23	24:BV:42:LEU:N	2.33	0.43
43:BW:45:HIS:HB2	43:BW:67:LYS:NZ	2.33	0.43
45:BY:23:LEU:HD22	45:BY:28:LEU:O	2.18	0.43
1:CA:1014:A:H5''	18:CS:13:HIS:HB2	2.00	0.43
1:CA:1426:G:O2'	1:CA:1427:C:H5'	2.19	0.43
1:CA:1455:G:O2'	1:CA:1456:A:H5'	2.18	0.43
1:CA:177:G:N3	1:CA:177:G:O4'	2.51	0.43
1:CA:302:G:H2'	1:CA:303:A:H8	1.83	0.43
1:CA:409:U:OP2	3:CD:21:LYS:HE2	2.19	0.43
20:CB:53:LEU:HD11	20:CB:216:VAL:HG12	1.99	0.43
3:CD:44:LYS:HZ2	3:CD:45:PRO:C	2.21	0.43
4:CE:87:VAL:CG1	4:CE:88:HIS:H	2.26	0.43
8:CI:123:ARG:HD3	8:CI:124:PRO:HD2	1.99	0.43
8:CI:8:THR:OG1	8:CI:84:ARG:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:625:U:H4'	15:CP:16:PHE:CZ	2.54	0.43
50:D3:41:ARG:HH11	50:D3:41:ARG:HG2	1.84	0.43
51:D4:2:LYS:HD3	51:D4:2:LYS:H	1.84	0.43
22:DA:53:A:H2'	22:DA:54:G:O4'	2.19	0.43
23:DB:1092:C:O2'	23:DB:1093:G:H5'	2.19	0.43
23:DB:1341:G:H5'	41:DT:61:LEU:CD2	2.44	0.43
23:DB:1373:A:H5''	23:DB:2213:U:O4	2.18	0.43
23:DB:1408:G:H2'	23:DB:1409:U:C6	2.54	0.43
23:DB:1547:C:H2'	23:DB:1548:A:C8	2.53	0.43
23:DB:1987:A:H2'	23:DB:1988:G:H8	1.82	0.43
23:DB:2135:A:C6	23:DB:2136:G:N7	2.87	0.43
23:DB:2260:C:O2'	23:DB:2261:C:H5'	2.18	0.43
23:DB:2581:G:N3	23:DB:2581:G:H2'	2.34	0.43
23:DB:2663:G:H2'	23:DB:2664:G:C8	2.54	0.43
23:DB:80:G:C6	23:DB:107:G:C6	3.07	0.43
23:DB:852:U:H2'	23:DB:853:C:C6	2.53	0.43
23:DB:948:C:O2'	23:DB:949:G:H5'	2.19	0.43
25:DC:45:ASN:ND2	25:DC:50:THR:OG1	2.51	0.43
28:DF:98:PHE:CA	28:DF:101:ARG:HE	2.23	0.43
28:DF:133:GLU:HB3	28:DF:134:GLN:H	1.50	0.43
29:DG:88:LEU:HG	29:DG:161:VAL:HB	2.01	0.43
29:DG:39:ALA:CA	29:DG:54:ARG:HB2	2.48	0.43
33:DL:74:THR:O	33:DL:75:ALA:C	2.57	0.43
34:DM:5:LYS:CE	34:DM:6:ARG:H	2.23	0.43
35:DN:100:CYS:SG	47:D0:43:THR:HG21	2.59	0.43
36:DO:35:ILE:CG1	36:DO:106:LEU:HD12	2.47	0.43
38:DQ:52:ARG:HG2	38:DQ:53:LYS:HE2	1.99	0.43
38:DQ:82:LEU:O	38:DQ:88:GLU:HG3	2.19	0.43
38:DQ:60:TRP:CD2	38:DQ:93:ILE:HA	2.54	0.43
46:DZ:12:ILE:HA	46:DZ:26:SER:HA	2.01	0.43
1:AA:1038:C:H2'	1:AA:1039:G:C8	2.54	0.43
1:AA:1489:G:H2'	1:AA:1490:U:C6	2.54	0.43
1:AA:223:A:H2'	1:AA:224:U:C6	2.53	0.43
1:AA:691:G:H1'	1:AA:696:A:H61	1.84	0.43
1:AA:967:C:OP1	1:AA:969:A:H8	2.00	0.43
20:AB:71:THR:HG23	20:AB:93:HIS:C	2.39	0.43
4:AE:56:PRO:O	4:AE:60:GLN:HG3	2.19	0.43
1:AA:1380:U:O4	6:AG:2:ARG:HG3	2.18	0.43
8:AI:94:ARG:CZ	8:AI:94:ARG:HB2	2.49	0.43
11:AL:64:SER:OG	11:AL:96:THR:HG23	2.18	0.43
12:AM:3:ILE:CA	12:AM:56:ARG:HG2	2.40	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AN:78:LEU:HB2	13:AN:83:VAL:HG22	1.99	0.43
17:AR:62:ARG:HD2	17:AR:69:TYR:CD2	2.54	0.43
21:AU:35:GLU:HB2	21:AU:37:TYR:CZ	2.54	0.43
23:BB:1306:C:H2'	23:BB:1307:A:H8	1.84	0.43
23:BB:1399:C:H2'	23:BB:1400:U:H6	1.84	0.43
23:BB:1549:A:H2'	23:BB:1550:C:H6	1.84	0.43
23:BB:1746:A:H2'	23:BB:1747:U:C6	2.53	0.43
23:BB:1857:G:H1'	23:BB:1885:A:H62	1.82	0.43
23:BB:1858:A:N6	23:BB:1884:G:O2'	2.51	0.43
23:BB:1945:G:C4	23:BB:1946:U:C5	3.07	0.43
23:BB:2100:G:C6	23:BB:2190:G:C6	3.07	0.43
23:BB:2233:U:H2'	23:BB:2234:G:H8	1.83	0.43
23:BB:2660:A:H2'	23:BB:2661:G:O4'	2.18	0.43
23:BB:2691:C:H2'	23:BB:2692:G:H8	1.83	0.43
23:BB:28:A:O2'	23:BB:29:U:H5'	2.18	0.43
23:BB:298:G:C2	23:BB:339:U:C5	3.06	0.43
25:BC:51:ARG:HB2	25:BC:51:ARG:HE	1.61	0.43
25:BC:66:PHE:HB3	25:BC:67:LYS:H	1.61	0.43
26:BD:1:MET:HG3	26:BD:205:PRO:HG2	2.01	0.43
28:BF:177:ARG:HH21	28:BF:178:LYS:HA	1.83	0.43
29:BG:8:VAL:O	29:BG:48:THR:HA	2.18	0.43
30:BH:30:LEU:HD22	30:BH:35:LYS:HE3	2.00	0.43
30:BH:5:LEU:CD1	30:BH:37:VAL:HG11	2.48	0.43
31:BJ:58:ASN:O	31:BJ:59:ALA:HB3	2.19	0.43
23:BB:872:U:HO2'	34:BM:64:TRP:HE1	1.62	0.43
35:BN:35:LYS:N	35:BN:112:TYR:HB3	2.33	0.43
35:BN:1:MET:HE3	35:BN:3:HIS:O	2.18	0.43
36:BO:1:MET:HG2	36:BO:2:ASP:N	2.32	0.43
36:BO:30:ARG:O	36:BO:31:THR:C	2.57	0.43
36:BO:50:ALA:HB1	36:BO:79:ALA:HB1	2.01	0.43
23:BB:1249:U:H4'	38:BQ:3:VAL:HG13	1.99	0.43
39:BR:51:VAL:O	39:BR:52:PRO:C	2.57	0.43
41:BT:11:LEU:HD23	41:BT:46:ALA:HB1	2.00	0.43
41:BT:3:ARG:HE	41:BT:8:LEU:HD23	1.83	0.43
43:BW:76:ARG:HG2	43:BW:77:LYS:H	1.81	0.43
43:BW:6:GLY:C	43:BW:8:SER:N	2.72	0.43
23:BB:188:G:OP1	46:BZ:11:GLU:HG2	2.18	0.43
1:CA:1013:G:H2'	1:CA:1015:G:OP2	2.18	0.43
1:CA:113:G:H21	1:CA:353:A:H8	1.66	0.43
1:CA:1298:U:H4'	1:CA:1299:A:C2	2.54	0.43
1:CA:1314:C:O2'	1:CA:1315:U:H5'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:186:C:H2'	1:CA:187:G:O4'	2.18	0.43
1:CA:709:U:H2'	1:CA:710:G:H8	1.83	0.43
1:CA:919:A:O2'	1:CA:920:U:H5'	2.19	0.43
20:CB:101:THR:HG22	20:CB:174:GLU:CD	2.39	0.43
2:CC:121:SER:O	2:CC:125:ARG:HG3	2.18	0.43
2:CC:54:ILE:HB	2:CC:67:ILE:CD1	2.49	0.43
3:CD:106:PHE:HB3	3:CD:154:VAL:HG12	2.01	0.43
6:CG:6:ILE:HG22	6:CG:7:GLY:N	2.34	0.43
7:CH:21:LYS:O	7:CH:62:LEU:HD22	2.19	0.43
8:CI:6:TYR:CG	8:CI:7:GLY:N	2.86	0.43
1:CA:1279:G:N2	9:CJ:45:ARG:HD2	2.33	0.43
9:CJ:59:LYS:C	9:CJ:61:ALA:H	2.21	0.43
12:CM:59:VAL:HG13	12:CM:60:ALA:N	2.33	0.43
14:CO:31:LEU:HA	14:CO:34:GLN:OE1	2.19	0.43
15:CP:13:LYS:O	15:CP:15:PRO:HD3	2.18	0.43
47:D0:16:ARG:C	47:D0:18:HIS:H	2.21	0.43
48:D1:9:LYS:HA	48:D1:24:LYS:HB3	2.00	0.43
22:DA:14:U:O2'	22:DA:107:G:H1'	2.19	0.43
22:DA:14:U:H3'	22:DA:15:A:C5'	2.48	0.43
23:DB:1199:U:O2'	38:DQ:2:ARG:HB2	2.18	0.43
23:DB:1270:C:H5''	23:DB:1271:G:O5'	2.19	0.43
23:DB:1281:G:H2'	23:DB:1282:U:O4'	2.18	0.43
23:DB:1559:U:H3'	23:DB:1560:G:C5'	2.47	0.43
23:DB:1599:U:C2	23:DB:1600:C:C5	3.06	0.43
23:DB:1707:G:H2'	23:DB:1708:C:C6	2.54	0.43
23:DB:1877:A:H2'	23:DB:1878:G:C8	2.53	0.43
23:DB:1858:A:N6	23:DB:1884:G:O2'	2.51	0.43
23:DB:1930:G:C2'	23:DB:1931:U:OP2	2.67	0.43
23:DB:1945:G:C4	23:DB:1946:U:C5	3.07	0.43
23:DB:2199:A:H3'	23:DB:2200:C:H6	1.82	0.43
23:DB:2601:C:C3'	23:DB:2602:A:H5''	2.47	0.43
23:DB:2678:C:O2'	23:DB:2679:A:H5'	2.18	0.43
23:DB:282:A:H2'	23:DB:283:G:C8	2.54	0.43
23:DB:480:A:C4'	42:DU:40:LEU:HD22	2.47	0.43
23:DB:839:U:H2'	23:DB:840:C:C6	2.53	0.43
25:DC:160:TYR:O	25:DC:176:ARG:HG2	2.19	0.43
26:DD:14:ILE:HG23	26:DD:19:GLY:CA	2.49	0.43
27:DE:136:GLN:HA	27:DE:139:LYS:HG2	2.01	0.43
28:DF:103:ILE:HD12	28:DF:104:THR:H	1.81	0.43
29:DG:29:ASN:HB3	29:DG:78:VAL:HA	2.00	0.43
32:DK:79:PHE:CD1	32:DK:79:PHE:N	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:14:LYS:HA	34:DM:72:PRO:HG3	2.00	0.43
35:DN:86:ARG:CZ	35:DN:117:ASP:HA	2.49	0.43
35:DN:80:PHE:O	35:DN:85:PRO:HD3	2.18	0.43
23:DB:585:G:N7	38:DQ:5:ARG:NH2	2.67	0.43
24:DV:44:HIS:O	24:DV:46:LYS:N	2.51	0.43
44:DX:39:GLN:O	44:DX:41:HIS:N	2.50	0.43
45:DY:5:LYS:HE2	45:DY:58:GLU:OE2	2.19	0.43
1:AA:1234:C:O2'	1:AA:1235:U:H5'	2.18	0.43
1:AA:1292:G:O2'	1:AA:1293:C:H5'	2.19	0.43
1:AA:408:A:OP1	3:AD:109:THR:HG21	2.19	0.43
1:AA:432:A:C5	1:AA:433:G:C8	3.07	0.43
1:AA:448:A:H2'	1:AA:449:G:C8	2.53	0.43
1:AA:374:A:H5''	1:AA:452:A:N1	2.33	0.43
1:AA:552:U:H5'	11:AL:82:ARG:HH11	1.83	0.43
1:AA:812:G:C2'	1:AA:812:G:N3	2.77	0.43
1:AA:995:C:O2	1:AA:1047:G:H5'	2.18	0.43
2:AC:112:ALA:HB2	2:AC:182:ASP:O	2.19	0.43
3:AD:84:ASN:OD1	4:AE:101:GLY:HA2	2.18	0.43
4:AE:109:ALA:O	4:AE:113:VAL:HG13	2.19	0.43
6:AG:110:ARG:HH12	6:AG:122:GLU:N	2.17	0.43
7:AH:95:MET:HB3	7:AH:99:GLY:H	1.84	0.43
8:AI:51:LEU:N	8:AI:51:LEU:HD12	2.34	0.43
9:AJ:93:ALA:HB1	9:AJ:96:VAL:CG2	2.49	0.43
10:AK:109:ILE:O	10:AK:110:THR:HG23	2.18	0.43
11:AL:41:PRO:HD3	11:AL:47:ALA:O	2.19	0.43
21:AU:44:ARG:HG3	21:AU:44:ARG:NH1	2.32	0.43
47:B0:30:ASP:CG	47:B0:31:LYS:H	2.22	0.43
23:BB:2479:U:OP1	51:B4:1:MET:SD	2.77	0.43
23:BB:1373:A:H4'	23:BB:2212:A:N3	2.34	0.43
23:BB:1426:G:H8	23:BB:1426:G:OP2	2.01	0.43
23:BB:2152:G:H2'	23:BB:2153:C:C6	2.54	0.43
23:BB:2247:A:O2'	23:BB:2248:C:H5'	2.19	0.43
23:BB:2272:U:C6	23:BB:2272:U:C3'	3.01	0.43
23:BB:2450:A:O2'	23:BB:2451:A:H5'	2.19	0.43
23:BB:2511:U:H2'	23:BB:2512:C:O4'	2.19	0.43
23:BB:2633:G:H2'	23:BB:2634:A:O4'	2.19	0.43
23:BB:2662:A:H2'	23:BB:2663:G:O4'	2.18	0.43
23:BB:281:C:H2'	23:BB:282:A:C8	2.54	0.43
23:BB:373:U:O2'	23:BB:374:A:H5'	2.19	0.43
23:BB:502:A:N6	23:BB:505:A:N1	2.66	0.43
25:BC:140:VAL:HG23	25:BC:141:HIS:N	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:152:GLN:HB3	25:BC:153:LEU:HD23	2.00	0.43
25:BC:164:VAL:HG12	25:BC:164:VAL:O	2.19	0.43
23:BB:2227:A:O2'	25:BC:258:SER:HA	2.19	0.43
25:BC:257:ARG:O	25:BC:261:ARG:HG3	2.18	0.43
26:BD:23:PRO:HA	26:BD:190:LYS:N	2.34	0.43
26:BD:31:ALA:HA	26:BD:51:THR:HG1	1.82	0.43
27:BE:158:PHE:C	27:BE:160:ALA:H	2.22	0.43
27:BE:59:PRO:HD2	27:BE:60:TRP:CE2	2.54	0.43
28:BF:55:ASP:OD2	28:BF:148:VAL:HG11	2.19	0.43
28:BF:90:LEU:N	28:BF:90:LEU:HD12	2.34	0.43
29:BG:121:THR:HG23	29:BG:133:LYS:O	2.18	0.43
29:BG:140:ILE:HD12	29:BG:141:GLY:N	2.34	0.43
52:BI:56:VAL:CG2	52:BI:68:PHE:HB2	2.49	0.43
31:BJ:57:LEU:HD12	31:BJ:128:ASN:HA	2.01	0.43
32:BK:107:LEU:C	32:BK:109:SER:H	2.22	0.43
32:BK:2:ILE:HG23	32:BK:33:ALA:O	2.19	0.43
34:BM:131:VAL:O	34:BM:132:THR:HB	2.18	0.43
34:BM:50:ARG:HH22	34:BM:54:THR:HG21	1.83	0.43
35:BN:29:VAL:HG13	35:BN:83:LEU:CD1	2.48	0.43
26:BD:8:LYS:CB	37:BP:5:LYS:HZ1	2.30	0.43
40:BS:28:LYS:HE3	40:BS:29:VAL:N	2.34	0.43
23:BB:492:A:H2	40:BS:7:HIS:HE1	1.67	0.43
43:BW:11:ASN:OD1	43:BW:12:GLY:N	2.52	0.43
1:CA:1190:G:OP2	2:CC:4:VAL:HB	2.19	0.43
1:CA:275:G:O5'	16:CQ:15:LYS:HG2	2.19	0.43
1:CA:432:A:C5	1:CA:433:G:C8	3.07	0.43
1:CA:607:A:H2'	1:CA:608:A:C8	2.52	0.43
1:CA:721:G:H4'	1:CA:722:G:O4'	2.19	0.43
5:CF:9:MET:HE1	17:CR:64:LEU:O	2.18	0.43
6:CG:100:MET:HA	6:CG:103:ILE:HD12	2.00	0.43
6:CG:144:ALA:C	6:CG:146:ALA:N	2.71	0.43
7:CH:68:LYS:HA	7:CH:68:LYS:HE3	2.01	0.43
9:CJ:85:ASP:O	9:CJ:86:ALA:C	2.57	0.43
1:CA:707:U:H4'	10:CK:21:HIS:CD2	2.54	0.43
13:CN:72:PHE:O	13:CN:73:LEU:HG	2.18	0.43
14:CO:25:GLU:CD	14:CO:76:ARG:HD3	2.38	0.43
15:CP:8:ARG:CZ	15:CP:15:PRO:HB3	2.48	0.43
18:CS:24:SER:HB2	18:CS:27:LYS:HZ1	1.83	0.43
18:CS:30:LEU:HB3	18:CS:48:ILE:HG23	2.00	0.43
47:D0:32:THR:HG21	47:D0:41:HIS:CD2	2.53	0.43
49:D2:19:ARG:HD3	49:D2:20:ALA:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:D4:18:LYS:O	51:D4:19:ARG:CB	2.67	0.43
23:DB:1062:G:H2'	23:DB:1063:G:H8	1.84	0.43
23:DB:1403:A:H2'	23:DB:1404:C:C6	2.54	0.43
23:DB:1426:G:H8	23:DB:1426:G:OP2	2.02	0.43
23:DB:1498:C:H2'	23:DB:1499:C:C6	2.53	0.43
23:DB:1923:U:H2'	23:DB:1924:C:H6	1.84	0.43
23:DB:2728:U:H2'	23:DB:2729:G:C8	2.53	0.43
23:DB:274:C:H2'	23:DB:275:C:C6	2.53	0.43
23:DB:2787:C:O2'	23:DB:2788:C:H5'	2.19	0.43
23:DB:2821:A:H2'	23:DB:2822:G:H8	1.82	0.43
23:DB:359:G:H2'	23:DB:360:U:H5'	2.00	0.43
23:DB:435:C:C2'	23:DB:436:C:H5'	2.49	0.43
23:DB:452:G:N2	23:DB:458:G:H1'	2.33	0.43
23:DB:565:C:O2'	23:DB:566:U:H5'	2.18	0.43
23:DB:693:A:H2'	23:DB:694:U:C6	2.53	0.43
23:DB:766:U:H2'	23:DB:767:U:H6	1.84	0.43
23:DB:909:A:H2'	23:DB:912:C:H5	1.83	0.43
25:DC:104:LEU:HD13	25:DC:156:SER:HB2	2.01	0.43
25:DC:155:ARG:HG2	25:DC:155:ARG:HH21	1.84	0.43
25:DC:21:PRO:HD2	25:DC:202:ARG:HH12	1.84	0.43
25:DC:90:ILE:O	25:DC:91:ALA:HB3	2.19	0.43
26:DD:128:ARG:O	26:DD:129:THR:HB	2.19	0.43
26:DD:173:GLN:HE21	26:DD:208:LYS:HB2	1.83	0.43
26:DD:42:ASN:O	26:DD:43:ASP:CB	2.64	0.43
26:DD:60:VAL:HG23	26:DD:63:PRO:CD	2.48	0.43
26:DD:47:ALA:H	26:DD:80:TRP:HB2	1.83	0.43
30:DH:69:ALA:O	30:DH:73:ASN:N	2.52	0.43
32:DK:4:GLU:O	32:DK:5:GLN:HB2	2.18	0.43
32:DK:71:ARG:O	32:DK:72:PRO:O	2.36	0.43
35:DN:34:ILE:O	35:DN:112:TYR:HD1	2.01	0.43
37:DP:111:GLU:O	37:DP:113:LEU:N	2.52	0.43
37:DP:27:VAL:HA	37:DP:86:LYS:HB3	2.01	0.43
38:DQ:7:VAL:HG13	38:DQ:8:ILE:CD1	2.42	0.43
40:DS:15:GLN:O	40:DS:19:LEU:HG	2.18	0.43
43:DW:33:GLY:HA3	43:DW:66:VAL:HG21	2.01	0.43
44:DX:31:GLN:NE2	44:DX:31:GLN:HA	2.29	0.43
44:DX:55:THR:O	44:DX:56:LEU:HB2	2.19	0.43
45:DY:26:LEU:HD21	45:DY:46:MET:HG2	2.00	0.43
46:DZ:3:LYS:CE	46:DZ:48:GLN:HB2	2.49	0.43
46:DZ:48:GLN:HB3	46:DZ:51:VAL:HG23	2.01	0.43
1:AA:1055:A:H4'	2:AC:160:GLU:HG3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:167:A:O2'	1:AA:168:G:H5'	2.19	0.43
1:AA:233:C:O2'	1:AA:234:C:H5'	2.19	0.43
1:AA:513:C:H2'	1:AA:514:C:C6	2.54	0.43
1:AA:664:G:H5''	17:AR:52:ARG:CZ	2.48	0.43
1:AA:709:U:H2'	1:AA:710:G:H8	1.84	0.43
20:AB:10:LYS:C	20:AB:12:GLY:H	2.22	0.43
20:AB:121:GLN:HB3	20:AB:121:GLN:HE21	1.52	0.43
2:AC:113:LYS:HZ3	2:AC:117:ASP:CG	2.22	0.43
2:AC:178:ARG:HG2	2:AC:178:ARG:O	2.19	0.43
2:AC:72:PRO:HG2	2:AC:73:GLY:H	1.84	0.43
3:AD:117:VAL:HG12	3:AD:130:ASN:HA	2.01	0.43
3:AD:28:ASP:O	3:AD:29:THR:OG1	2.26	0.43
3:AD:81:LEU:HD11	3:AD:92:LEU:HD21	2.00	0.43
7:AH:17:GLN:OE1	7:AH:69:ALA:HB1	2.19	0.43
7:AH:9:MET:HE3	7:AH:32:LYS:HB3	2.01	0.43
9:AJ:89:ARG:O	9:AJ:90:LEU:HD23	2.19	0.43
11:AL:56:LEU:HD11	11:AL:81:ILE:CD1	2.41	0.43
13:AN:11:LYS:NZ	13:AN:11:LYS:HA	2.34	0.43
14:AO:2:LEU:HD22	14:AO:7:THR:HG23	2.00	0.43
16:AQ:60:ILE:HB	16:AQ:72:TRP:CE3	2.54	0.43
21:AU:16:ARG:C	21:AU:18:PHE:H	2.21	0.43
23:BB:100:U:OP1	23:BB:100:U:H2'	2.18	0.43
23:BB:1076:C:O2'	23:BB:1077:A:H5'	2.19	0.43
23:BB:1081:U:O2'	23:BB:1082:U:H5'	2.19	0.43
23:BB:1281:G:H2'	23:BB:1282:U:O4'	2.18	0.43
23:BB:1704:C:H2'	23:BB:1705:A:C8	2.54	0.43
23:BB:1707:G:H2'	23:BB:1708:C:C6	2.54	0.43
23:BB:357:C:H2'	23:BB:358:U:C6	2.54	0.43
23:BB:36:G:H4'	23:BB:451:U:N3	2.34	0.43
23:BB:756:A:H2'	23:BB:757:G:O4'	2.19	0.43
25:BC:153:LEU:O	25:BC:153:LEU:HG	2.19	0.43
25:BC:75:ALA:HB1	25:BC:93:VAL:HG13	2.01	0.43
27:BE:153:LEU:HD11	27:BE:201:ALA:HA	2.00	0.43
28:BF:2:LYS:CA	28:BF:2:LYS:HE2	2.48	0.43
29:BG:137:LYS:HA	29:BG:140:ILE:HG13	2.00	0.43
30:BH:89:LYS:NZ	30:BH:123:ARG:HD3	2.34	0.43
30:BH:28:ASN:HA	30:BH:32:PRO:HG3	2.00	0.43
27:BE:25:GLU:OE2	33:BL:16:GLY:HA2	2.19	0.43
33:BL:60:ARG:HD3	50:B3:11:LYS:CE	2.49	0.43
34:BM:36:VAL:HG21	34:BM:125:PRO:HD3	2.00	0.43
34:BM:43:ALA:HB2	34:BM:91:TYR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:57:THR:CB	35:BN:61:ALA:HB3	2.49	0.43
37:BP:28:LYS:C	37:BP:44:GLY:HA3	2.39	0.43
38:BQ:102:LYS:H	38:BQ:102:LYS:CD	2.29	0.43
38:BQ:33:VAL:O	38:BQ:36:GLN:N	2.50	0.43
38:BQ:35:PHE:HA	38:BQ:39:ILE:HD12	2.00	0.43
42:BU:70:ALA:O	42:BU:71:ILE:C	2.56	0.43
42:BU:76:THR:O	42:BU:77:GLY:C	2.56	0.43
42:BU:78:LYS:CG	42:BU:96:LYS:HE3	2.48	0.43
46:BZ:3:LYS:HA	46:BZ:47:LYS:HA	2.00	0.43
46:BZ:4:ASP:O	46:BZ:5:ILE:C	2.57	0.43
1:CA:1067:A:H3'	1:CA:1094:G:OP1	2.19	0.43
1:CA:513:C:H2'	1:CA:514:C:C6	2.53	0.43
1:CA:730:G:O6	14:CO:50:HIS:NE2	2.49	0.43
1:CA:822:U:O2'	1:CA:823:C:H5'	2.19	0.43
1:CA:979:C:H41	1:CA:1360:A:H62	1.67	0.43
20:CB:216:VAL:HG23	20:CB:217:ALA:N	2.34	0.43
3:CD:44:LYS:NZ	3:CD:46:ARG:HA	2.34	0.43
10:CK:63:GLN:HG3	10:CK:98:ALA:HB2	2.01	0.43
12:CM:37:GLY:O	12:CM:38:ILE:HD13	2.18	0.43
13:CN:73:LEU:O	13:CN:74:ARG:C	2.57	0.43
47:D0:46:GLY:O	47:D0:47:TYR:HB3	2.18	0.43
50:D3:39:ARG:HD3	50:D3:39:ARG:HA	1.76	0.43
22:DA:32:U:C2	22:DA:33:G:C8	3.07	0.43
23:DB:1704:C:H2'	23:DB:1705:A:C8	2.54	0.43
23:DB:1829:A:H3'	23:DB:1830:C:H6	1.84	0.43
1:CA:1409:C:O2	23:DB:1913:A:N6	2.52	0.43
23:DB:2157:G:N3	23:DB:2157:G:C2'	2.79	0.43
23:DB:2361:G:O2'	23:DB:2362:C:H5'	2.18	0.43
23:DB:2466:C:O2'	23:DB:2467:C:H5'	2.17	0.43
23:DB:2564:A:OP1	23:DB:2648:G:H4'	2.18	0.43
23:DB:593:U:H2'	23:DB:594:U:H6	1.81	0.43
23:DB:878:A:H1'	23:DB:899:A:H62	1.83	0.43
26:DD:33:ARG:C	26:DD:35:THR:H	2.22	0.43
27:DE:176:ASP:OD1	27:DE:177:PRO:HD2	2.19	0.43
29:DG:8:VAL:CG1	29:DG:49:LEU:HD12	2.48	0.43
52:DI:54:ILE:HD11	52:DI:71:LYS:O	2.19	0.43
32:DK:10:VAL:HG21	32:DK:17:ARG:N	2.34	0.43
33:DL:107:PHE:HE2	33:DL:126:ARG:HB2	1.84	0.43
33:DL:14:LYS:O	33:DL:15:ALA:C	2.57	0.43
23:DB:598:U:C5'	33:DL:21:ARG:HB2	2.45	0.43
23:DB:631:A:O2'	33:DL:66:PHE:CD1	2.70	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:73:ILE:CD1	34:DM:92:TRP:HE3	2.32	0.43
38:DQ:25:GLY:O	38:DQ:29:ARG:HG3	2.18	0.43
39:DR:68:ARG:H	39:DR:97:LYS:HG3	1.84	0.43
39:DR:18:GLN:HB3	39:DR:99:THR:HB	2.01	0.43
42:DU:46:LYS:HG2	42:DU:53:GLN:HB2	2.01	0.43
44:DX:17:GLU:CA	44:DX:21:LEU:HB2	2.49	0.43
45:DY:35:VAL:HG12	45:DY:36:GLU:H	1.83	0.43
46:DZ:49:ARG:O	46:DZ:52:ALA:N	2.50	0.43
1:AA:1263:C:H2'	1:AA:1264:U:H6	1.84	0.43
1:AA:1322:C:O4'	1:AA:1322:C:O2	2.37	0.43
1:AA:155:A:H2'	1:AA:156:C:O4'	2.19	0.43
1:AA:25:C:H2'	1:AA:26:A:C8	2.54	0.43
1:AA:371:A:H2'	1:AA:372:C:C6	2.54	0.43
1:AA:401:C:O2'	1:AA:402:G:H5'	2.18	0.43
1:AA:484:G:O2'	1:AA:485:U:OP2	2.29	0.43
1:AA:529:G:H22	11:AL:47:ALA:HB3	1.84	0.43
1:AA:626:G:H2'	1:AA:627:G:H8	1.83	0.43
1:AA:659:U:O2'	1:AA:660:C:H5'	2.18	0.43
1:AA:721:G:H4'	1:AA:722:G:O4'	2.19	0.43
1:AA:861:G:H2'	1:AA:862:C:H6	1.84	0.43
20:AB:217:ALA:O	20:AB:221:ARG:HG3	2.18	0.43
20:AB:35:ASN:O	20:AB:36:LYS:HB2	2.19	0.43
5:AF:7:VAL:HG23	5:AF:60:VAL:O	2.18	0.43
8:AI:82:ILE:HG22	8:AI:86:LEU:HD21	2.01	0.43
9:AJ:57:VAL:HG13	9:AJ:58:ASN:CG	2.39	0.43
9:AJ:57:VAL:HG13	9:AJ:58:ASN:N	2.33	0.43
9:AJ:7:ARG:HG2	9:AJ:75:ASP:OD2	2.18	0.43
16:AQ:10:ARG:HH22	16:AQ:58:VAL:HG13	1.84	0.43
47:B0:4:GLN:O	47:B0:5:ASN:C	2.57	0.43
49:B2:3:ARG:HB3	49:B2:5:PHE:HE1	1.84	0.43
50:B3:22:LYS:HD2	50:B3:46:LYS:HA	1.99	0.43
23:BB:1359:A:H2'	23:BB:1360:G:O4'	2.17	0.43
23:BB:1634:A:H3'	23:BB:1635:A:C5'	2.48	0.43
23:BB:1773:A:H2'	23:BB:1774:C:O4'	2.19	0.43
23:BB:2511:U:O5'	23:BB:2511:U:H6	2.02	0.43
23:BB:2663:G:H2'	23:BB:2664:G:C8	2.54	0.43
23:BB:2671:G:O2'	23:BB:2672:U:H5'	2.19	0.43
23:BB:2811:G:H2'	23:BB:2812:G:H8	1.83	0.43
23:BB:285:G:C2'	23:BB:286:U:H5'	2.49	0.43
23:BB:511:U:H6	23:BB:511:U:O5'	2.01	0.43
23:BB:832:U:OP1	33:BL:45:GLY:N	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:848:C:H3'	23:BB:849:A:H8	1.83	0.43
23:BB:971:G:H2'	23:BB:972:A:O4'	2.19	0.43
23:BB:975:A:H1'	23:BB:990:A:C2	2.54	0.43
25:BC:156:SER:O	25:BC:157:ALA:C	2.58	0.43
25:BC:225:ASN:HD22	25:BC:225:ASN:HA	1.52	0.43
25:BC:247:TRP:HD1	25:BC:247:TRP:O	2.02	0.43
27:BE:144:GLU:HG3	27:BE:145:ASP:N	2.34	0.43
27:BE:189:THR:O	27:BE:190:ALA:HB3	2.19	0.43
29:BG:140:ILE:HD12	29:BG:140:ILE:C	2.39	0.43
52:BI:75:ALA:CB	52:BI:131:THR:HG21	2.49	0.43
31:BJ:80:HIS:HB2	31:BJ:84:ILE:HD11	1.99	0.43
32:BK:108:ARG:HH11	32:BK:108:ARG:HG3	1.83	0.43
23:BB:1244:A:H4'	33:BL:15:ALA:HB1	2.00	0.43
33:BL:61:LEU:CB	33:BL:62:PRO:CD	2.92	0.43
35:BN:19:ALA:C	35:BN:21:PHE:N	2.72	0.43
35:BN:99:LYS:CB	35:BN:99:LYS:NZ	2.79	0.43
36:BO:67:ASN:HD22	36:BO:67:ASN:HA	1.65	0.43
37:BP:51:ASN:HA	37:BP:60:VAL:HB	2.00	0.43
40:BS:29:VAL:H	40:BS:71:VAL:HG21	1.84	0.43
40:BS:33:LEU:O	40:BS:35:ILE:N	2.50	0.43
23:BB:2330:G:O2'	43:BW:36:ILE:HD11	2.18	0.43
44:BX:18:LEU:HD23	44:BX:18:LEU:N	2.34	0.43
1:CA:113:G:C1'	1:CA:354:G:H5'	2.49	0.43
1:CA:815:A:N6	1:CA:1509:C:H1'	2.32	0.43
1:CA:179:A:H2'	1:CA:180:U:C6	2.54	0.43
1:CA:408:A:H3'	1:CA:409:U:C6	2.47	0.43
1:CA:453:G:C6	1:CA:454:G:C6	3.06	0.43
1:CA:635:A:H2'	1:CA:636:U:C6	2.53	0.43
1:CA:814:A:H5'	1:CA:1511:G:C4'	2.42	0.43
20:CB:111:LYS:O	20:CB:114:LYS:HB3	2.19	0.43
2:CC:166:TRP:HE1	2:CC:168:ARG:CB	2.32	0.43
3:CD:121:ALA:C	3:CD:145:ARG:HG3	2.39	0.43
3:CD:21:LYS:C	3:CD:23:GLY:H	2.22	0.43
6:CG:29:LEU:CD2	6:CG:104:VAL:HG13	2.49	0.43
7:CH:50:VAL:HG23	7:CH:57:GLU:O	2.19	0.43
11:CL:109:ARG:C	11:CL:110:LYS:HD2	2.39	0.43
1:CA:1309:G:H5'	12:CM:76:ILE:HG12	2.00	0.43
13:CN:26:LEU:O	13:CN:44:VAL:HG11	2.19	0.43
16:CQ:35:LYS:O	16:CQ:37:ILE:HG23	2.19	0.43
51:D4:19:ARG:HB3	51:D4:20:ASP:H	1.47	0.43
22:DA:111:U:O2'	22:DA:112:G:H5'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1841:U:C2	23:DB:1842:G:C8	3.07	0.43
23:DB:1891:G:H2'	23:DB:1892:C:H6	1.84	0.43
23:DB:2256:G:H2'	23:DB:2257:U:H6	1.84	0.43
23:DB:2341:G:H2'	23:DB:2342:C:C6	2.54	0.43
23:DB:2405:G:H5'	33:DL:70:LYS:CG	2.49	0.43
23:DB:2450:A:O2'	23:DB:2451:A:H5'	2.19	0.43
23:DB:2782:G:O2'	23:DB:2783:U:H5'	2.19	0.43
23:DB:539:G:H2'	23:DB:540:C:C6	2.54	0.43
23:DB:630:G:N1	33:DL:69:ARG:NH1	2.66	0.43
23:DB:981:A:N1	23:DB:2027:G:O2'	2.43	0.43
25:DC:106:PRO:HB2	25:DC:107:LYS:H	1.68	0.43
25:DC:235:GLU:CG	25:DC:236:GLY:H	2.30	0.43
26:DD:172:VAL:HB	26:DD:175:LEU:HD11	2.00	0.43
26:DD:25:THR:HG22	26:DD:188:LEU:CD1	2.49	0.43
26:DD:4:LEU:HB3	26:DD:202:ILE:HA	2.01	0.43
30:DH:4:ILE:CD1	30:DH:37:VAL:HG13	2.49	0.43
33:DL:79:LEU:H	33:DL:113:ALA:CB	2.31	0.43
34:DM:133:LYS:CD	34:DM:134:THR:H	2.30	0.43
35:DN:21:PHE:HA	35:DN:24:MET:HB3	2.01	0.43
35:DN:4:ARG:HB3	35:DN:4:ARG:CZ	2.49	0.43
37:DP:20:ARG:HH21	37:DP:20:ARG:HG2	1.84	0.43
39:DR:38:VAL:O	39:DR:41:ILE:HG22	2.18	0.43
39:DR:76:LYS:HE3	39:DR:90:ARG:HD3	2.01	0.43
41:DT:47:VAL:CG2	41:DT:53:VAL:HG21	2.42	0.43
43:DW:17:ALA:O	43:DW:18:LYS:CG	2.66	0.43
46:DZ:3:LYS:HE2	46:DZ:48:GLN:HB2	2.00	0.43
1:AA:1035:A:C2'	1:AA:1036:A:H8	2.27	0.42
1:AA:1072:G:H2'	1:AA:1073:U:O4'	2.19	0.42
1:AA:113:G:C1'	1:AA:354:G:H5'	2.49	0.42
1:AA:179:A:H2'	1:AA:180:U:H6	1.83	0.42
1:AA:208:U:H2'	1:AA:210:C:C5	2.54	0.42
1:AA:332:G:P	19:AT:2:ASN:HB3	2.59	0.42
1:AA:349:A:H2'	1:AA:350:G:C8	2.54	0.42
1:AA:538:G:H2'	1:AA:539:A:C8	2.54	0.42
1:AA:997:U:H2'	1:AA:998:C:H6	1.82	0.42
20:AB:21:TYR:O	20:AB:22:TRP:O	2.36	0.42
20:AB:37:VAL:CG2	20:AB:38:HIS:N	2.81	0.42
3:AD:138:PRO:HA	3:AD:181:PHE:HD2	1.83	0.42
3:AD:201:GLU:CG	4:AE:103:GLY:HA3	2.49	0.42
6:AG:117:LEU:C	6:AG:117:LEU:HD13	2.39	0.42
8:AI:115:VAL:HG23	9:AJ:62:ARG:HD2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:106:VAL:CG2	11:AL:109:ARG:HG3	2.48	0.42
11:AL:58:ASN:H	11:AL:58:ASN:HD22	1.66	0.42
14:AO:71:ARG:HH11	14:AO:71:ARG:CB	2.32	0.42
18:AS:29:PRO:HB3	18:AS:47:THR:HB	2.00	0.42
19:AT:4:LYS:HD2	19:AT:5:SER:H	1.82	0.42
19:AT:50:PHE:CE2	19:AT:75:LYS:HA	2.53	0.42
47:B0:53:VAL:HG12	47:B0:54:ILE:N	2.34	0.42
48:B1:47:ILE:HG13	48:B1:48:TYR:N	2.34	0.42
48:B1:53:ILE:CD1	48:B1:54:LYS:H	2.32	0.42
23:BB:241:A:H4'	50:B3:63:TYR:OH	2.18	0.42
23:BB:1062:G:H2'	23:BB:1063:G:H8	1.84	0.42
23:BB:1079:C:C2	23:BB:1080:A:C8	3.07	0.42
23:BB:1314:C:O2	23:BB:1314:C:H2'	2.18	0.42
23:BB:1526:C:O2'	23:BB:1527:G:H5'	2.19	0.42
23:BB:1917:U:H2'	23:BB:1918:A:C5'	2.48	0.42
23:BB:2290:G:H2'	23:BB:2291:U:C6	2.53	0.42
23:BB:402:A:C2'	23:BB:403:U:H5'	2.49	0.42
25:BC:198:GLU:HA	25:BC:201:LEU:HB2	2.01	0.42
25:BC:78:GLU:CB	25:BC:92:LEU:HD23	2.49	0.42
26:BD:108:ASP:CB	26:BD:173:GLN:HA	2.48	0.42
26:BD:130:GLN:O	26:BD:140:HIS:ND1	2.51	0.42
23:BB:2052:A:C4	26:BD:155:VAL:HG23	2.54	0.42
27:BE:128:ALA:CA	27:BE:157:LEU:HD22	2.48	0.42
28:BF:15:LEU:O	28:BF:18:GLU:O	2.37	0.42
28:BF:67:THR:OG1	28:BF:85:GLY:HA3	2.19	0.42
29:BG:5:LYS:O	29:BG:68:ARG:NH1	2.52	0.42
31:BJ:102:GLU:CG	31:BJ:124:VAL:HG11	2.49	0.42
31:BJ:128:ASN:CG	31:BJ:129:GLU:N	2.72	0.42
23:BB:995:C:C4	31:BJ:2:LYS:HB3	2.54	0.42
31:BJ:7:LYS:HD2	31:BJ:49:ASP:CB	2.44	0.42
32:BK:71:ARG:HD3	32:BK:71:ARG:HA	1.87	0.42
33:BL:38:GLN:HA	33:BL:41:ARG:NH1	2.34	0.42
33:BL:63:LYS:C	50:B3:24:LYS:HZ2	2.23	0.42
34:BM:134:THR:O	34:BM:135:VAL:HG13	2.18	0.42
23:BB:2333:A:O2'	36:BO:7:ARG:HD3	2.19	0.42
41:BT:47:VAL:HB	41:BT:55:VAL:HG21	2.00	0.42
41:BT:50:LEU:O	41:BT:51:PHE:HB2	2.18	0.42
23:BB:2270:A:C5'	43:BW:18:LYS:HB3	2.49	0.42
43:BW:35:ILE:CG2	43:BW:36:ILE:H	2.14	0.42
1:CA:1264:U:H2'	1:CA:1265:C:O4'	2.19	0.42
1:CA:975:A:H1'	1:CA:1358:U:O2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1521:C:H2'	1:CA:1522:U:H6	1.83	0.42
1:CA:179:A:H2'	1:CA:180:U:H6	1.84	0.42
1:CA:406:G:H21	3:CD:115:GLN:HE22	1.65	0.42
1:CA:697:U:C2'	1:CA:698:G:H5'	2.49	0.42
1:CA:69:G:N2	1:CA:71:A:H62	2.16	0.42
1:CA:861:G:H2'	1:CA:862:C:H6	1.84	0.42
1:CA:994:A:H2'	1:CA:994:A:N3	2.34	0.42
20:CB:121:GLN:NE2	20:CB:122:ASP:N	2.67	0.42
2:CC:168:ARG:HE	2:CC:169:GLU:N	2.17	0.42
3:CD:170:LEU:CA	3:CD:182:LYS:HG3	2.48	0.42
5:CF:3:HIS:CB	5:CF:92:THR:HA	2.31	0.42
8:CI:21:LYS:O	8:CI:61:ASP:N	2.51	0.42
9:CJ:40:ILE:HG12	9:CJ:74:VAL:H	1.84	0.42
9:CJ:8:ILE:HG12	9:CJ:75:ASP:HA	2.00	0.42
9:CJ:9:ARG:HH11	9:CJ:9:ARG:HA	1.84	0.42
11:CL:17:LYS:HZ3	11:CL:17:LYS:HB2	1.82	0.42
13:CN:50:LEU:HD12	13:CN:51:PRO:CD	2.46	0.42
13:CN:13:VAL:HA	13:CN:59:GLN:HE21	1.84	0.42
14:CO:80:LEU:C	14:CO:80:LEU:HD23	2.39	0.42
17:CR:20:ILE:HG23	17:CR:20:ILE:O	2.19	0.42
40:DS:23:LEU:CB	47:D0:21:LEU:HD13	2.37	0.42
48:D1:24:LYS:O	48:D1:24:LYS:HD3	2.19	0.42
22:DA:60:C:H2'	22:DA:61:G:H8	1.83	0.42
23:DB:1513:U:O2'	23:DB:1514:G:H5'	2.19	0.42
23:DB:153:U:O5'	23:DB:153:U:H6	2.02	0.42
23:DB:1770:G:O2'	23:DB:1771:C:H5'	2.19	0.42
23:DB:1803:A:H2	23:DB:1823:G:O4'	2.02	0.42
23:DB:758:C:O2	23:DB:1981:A:H2	2.01	0.42
23:DB:1998:A:H2'	23:DB:1999:C:H6	1.84	0.42
23:DB:2094:A:H5''	30:DH:22:LYS:HG3	2.01	0.42
23:DB:2076:U:OP2	23:DB:2238:G:N2	2.48	0.42
23:DB:2336:A:HO2'	23:DB:2337:G:P	2.41	0.42
23:DB:2385:C:H2'	23:DB:2386:A:C8	2.53	0.42
23:DB:2428:G:H5''	23:DB:2429:G:O5'	2.18	0.42
23:DB:2511:U:H2'	23:DB:2512:C:O4'	2.19	0.42
23:DB:2513:A:H2'	23:DB:2514:U:C6	2.54	0.42
23:DB:2671:G:H2'	23:DB:2672:U:H6	1.80	0.42
23:DB:2811:G:H2'	23:DB:2812:G:C8	2.54	0.42
23:DB:2813:A:O2'	23:DB:2814:A:H5'	2.18	0.42
23:DB:409:G:O2'	23:DB:410:G:H5'	2.19	0.42
23:DB:818:G:H4'	23:DB:838:C:O3'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:172:THR:HB	25:DC:173:LEU:H	1.56	0.42
26:DD:109:VAL:HA	26:DD:203:VAL:HA	2.01	0.42
26:DD:113:SER:O	26:DD:167:ASN:HB2	2.19	0.42
26:DD:11:MET:O	26:DD:22:ILE:HD12	2.19	0.42
28:DF:41:GLU:OE1	28:DF:41:GLU:N	2.52	0.42
30:DH:12:LEU:HD22	30:DH:19:VAL:CG1	2.49	0.42
30:DH:77:THR:HG22	30:DH:143:ILE:HB	2.00	0.42
52:DI:49:GLU:HB2	52:DI:52:LEU:HD12	2.00	0.42
31:DJ:62:VAL:HG11	31:DJ:101:ILE:HD11	2.00	0.42
33:DL:25:SER:C	33:DL:27:LEU:N	2.72	0.42
36:DO:71:ALA:O	36:DO:74:VAL:HG22	2.19	0.42
23:DB:2294:G:P	36:DO:9:ARG:NE	2.92	0.42
37:DP:32:VAL:O	37:DP:34:GLY:N	2.51	0.42
38:DQ:32:ARG:HG2	38:DQ:32:ARG:NH1	2.32	0.42
39:DR:67:GLY:H	39:DR:98:ILE:N	2.17	0.42
39:DR:69:GLY:HA3	39:DR:96:VAL:HG13	2.00	0.42
41:DT:34:VAL:O	41:DT:81:LYS:HB3	2.19	0.42
43:DW:36:ILE:HG13	43:DW:37:VAL:N	2.34	0.42
45:DY:5:LYS:O	45:DY:56:VAL:HG12	2.19	0.42
45:DY:5:LYS:NZ	45:DY:5:LYS:HB2	2.34	0.42
1:AA:1082:A:H2'	1:AA:1083:U:C6	2.53	0.42
1:AA:113:G:O4'	1:AA:354:G:H4'	2.19	0.42
1:AA:425:G:O2'	1:AA:426:U:H5'	2.18	0.42
1:AA:523:A:H61	11:AL:88:ASP:HB2	1.84	0.42
1:AA:644:U:O2'	1:AA:645:G:H5'	2.20	0.42
1:AA:697:U:C2'	1:AA:698:G:H5'	2.49	0.42
1:AA:738:C:H2'	1:AA:739:C:C6	2.55	0.42
1:AA:822:U:O2'	1:AA:823:C:H5'	2.19	0.42
1:AA:93:U:H2'	1:AA:95:C:H5	1.83	0.42
3:AD:29:THR:O	3:AD:30:LYS:HB2	2.19	0.42
7:AH:98:LEU:N	7:AH:98:LEU:HD12	2.34	0.42
8:AI:49:GLN:HB3	8:AI:49:GLN:HE21	1.61	0.42
12:AM:79:LEU:HD22	12:AM:86:ARG:NH2	2.33	0.42
15:AP:52:LEU:HG	15:AP:75:ILE:CG1	2.48	0.42
16:AQ:28:VAL:HG13	16:AQ:28:VAL:O	2.19	0.42
19:AT:60:GLN:HB2	19:AT:65:LEU:HD12	2.01	0.42
19:AT:70:LYS:HA	19:AT:73:ARG:CZ	2.48	0.42
47:B0:53:VAL:HG12	47:B0:54:ILE:H	1.84	0.42
23:BB:2082:A:H2'	23:BB:2083:G:O4'	2.19	0.42
23:BB:2570:G:O2'	23:BB:2571:U:H5'	2.19	0.42
23:BB:2637:U:O2'	23:BB:2638:G:H5'	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:766:U:H2'	23:BB:767:U:H6	1.84	0.42
23:BB:822:G:H2'	23:BB:823:C:H6	1.84	0.42
25:BC:27:LYS:HA	25:BC:79:ARG:NH2	2.34	0.42
25:BC:84:PRO:O	25:BC:86:ARG:N	2.48	0.42
26:BD:142:VAL:C	26:BD:144:GLY:H	2.21	0.42
27:BE:135:ALA:O	27:BE:139:LYS:HD2	2.19	0.42
28:BF:177:ARG:HE	28:BF:178:LYS:HA	1.81	0.42
30:BH:31:VAL:HB	30:BH:32:PRO:HD3	2.01	0.42
30:BH:80:ILE:HD11	30:BH:102:ALA:HA	2.01	0.42
52:BI:116:MET:CE	52:BI:127:SER:HB2	2.49	0.42
33:BL:127:VAL:HB	33:BL:128:THR:H	1.51	0.42
33:BL:96:LYS:N	33:BL:96:LYS:HD3	2.34	0.42
34:BM:36:VAL:HG11	34:BM:125:PRO:HD2	2.00	0.42
34:BM:25:ASP:OD2	34:BM:66:ARG:NH2	2.52	0.42
35:BN:19:ALA:C	35:BN:21:PHE:H	2.22	0.42
37:BP:61:ARG:O	37:BP:63:ILE:N	2.52	0.42
37:BP:7:LEU:C	37:BP:7:LEU:HD13	2.39	0.42
39:BR:39:LEU:O	39:BR:39:LEU:HD23	2.18	0.42
39:BR:53:PHE:CD2	39:BR:55:ASP:N	2.82	0.42
41:BT:3:ARG:HG2	41:BT:7:LEU:HD23	2.01	0.42
41:BT:69:ARG:H	41:BT:75:GLY:CA	2.32	0.42
42:BU:13:LEU:H	42:BU:68:ASN:HD21	1.66	0.42
43:BW:9:THR:H	43:BW:10:ARG:HD3	1.83	0.42
46:BZ:58:ASP:OD2	46:BZ:63:ARG:NH1	2.52	0.42
1:CA:1065:U:O2'	1:CA:1066:C:OP2	2.30	0.42
1:CA:112:G:O2'	1:CA:113:G:H5'	2.19	0.42
1:CA:1493:A:C2'	1:CA:1494:G:OP2	2.67	0.42
1:CA:1531:A:H2'	1:CA:1532:U:O4'	2.19	0.42
1:CA:79:G:H2'	1:CA:80:A:C8	2.53	0.42
20:CB:128:LEU:HD21	20:CB:132:GLU:HB3	2.01	0.42
20:CB:113:LEU:HD11	20:CB:144:GLU:HA	2.01	0.42
2:CC:137:VAL:HG13	2:CC:148:ILE:CG2	2.49	0.42
2:CC:64:ARG:O	2:CC:65:VAL:HB	2.19	0.42
4:CE:89:THR:O	4:CE:91:SER:N	2.52	0.42
5:CF:11:HIS:CD2	5:CF:13:ASP:HB3	2.54	0.42
6:CG:46:LEU:O	6:CG:50:ALA:HB2	2.19	0.42
9:CJ:84:VAL:O	9:CJ:88:MET:HB3	2.19	0.42
9:CJ:91:ASP:C	9:CJ:92:LEU:HD23	2.40	0.42
10:CK:86:LYS:HG3	10:CK:113:THR:HA	2.00	0.42
15:CP:48:GLU:HG2	15:CP:49:GLY:N	2.34	0.42
15:CP:5:ARG:O	15:CP:19:VAL:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:CQ:75:VAL:HG23	16:CQ:76:ARG:N	2.33	0.42
21:CU:19:LYS:HG3	21:CU:20:ARG:HE	1.83	0.42
47:D0:15:ARG:HB3	47:D0:15:ARG:CZ	2.49	0.42
51:D4:24:ARG:NE	51:D4:37:GLN:HA	2.33	0.42
23:DB:1068:G:C6	23:DB:1069:A:N6	2.87	0.42
23:DB:1260:A:O2'	23:DB:1261:C:H5'	2.19	0.42
23:DB:1818:U:C3'	25:DC:155:ARG:HB2	2.49	0.42
23:DB:2087:G:H2'	23:DB:2088:A:H8	1.85	0.42
23:DB:2508:G:O3'	23:DB:2555:U:H5'	2.18	0.42
23:DB:2649:C:O2'	23:DB:2650:U:H5'	2.19	0.42
23:DB:2866:U:O3'	23:DB:2867:G:H4'	2.19	0.42
23:DB:329:G:H1'	23:DB:477:A:H1'	2.02	0.42
23:DB:348:A:H2'	23:DB:349:U:O4'	2.18	0.42
23:DB:543:G:C6	23:DB:544:C:H1'	2.54	0.42
23:DB:679:C:O2'	23:DB:680:C:H5'	2.19	0.42
23:DB:96:C:H2'	23:DB:97:C:H6	1.84	0.42
25:DC:107:LYS:HD2	25:DC:196:ASN:HD21	1.84	0.42
25:DC:62:ARG:HH22	25:DC:86:ARG:HH21	1.66	0.42
26:DD:154:LYS:O	26:DD:156:PHE:N	2.51	0.42
27:DE:146:VAL:O	27:DE:148:ILE:N	2.51	0.42
27:DE:123:LYS:HD2	27:DE:157:LEU:CD1	2.48	0.42
27:DE:49:ARG:C	27:DE:51:GLU:N	2.73	0.42
29:DG:136:ASP:O	29:DG:140:ILE:HG12	2.19	0.42
29:DG:154:GLU:OE2	29:DG:157:LYS:HB2	2.18	0.42
29:DG:23:ILE:HD13	29:DG:71:LEU:HD22	2.01	0.42
30:DH:12:LEU:C	30:DH:12:LEU:HD23	2.39	0.42
30:DH:93:SER:C	30:DH:94:ILE:HD12	2.39	0.42
33:DL:77:ILE:CD1	33:DL:109:LYS:HB2	2.46	0.42
23:DB:631:A:O2'	33:DL:66:PHE:HD1	1.97	0.42
34:DM:41:LEU:HD21	34:DM:46:ILE:HD11	2.01	0.42
34:DM:43:ALA:HB3	34:DM:91:TYR:CD2	2.55	0.42
36:DO:86:GLY:C	36:DO:88:LYS:N	2.73	0.42
37:DP:92:ARG:HD3	37:DP:110:LYS:O	2.19	0.42
38:DQ:39:ILE:C	38:DQ:39:ILE:HD12	2.40	0.42
38:DQ:90:ASP:CB	39:DR:42:ALA:HB3	2.49	0.42
44:DX:46:VAL:O	44:DX:49:ASP:HB3	2.18	0.42
1:AA:1250:A:H5'	8:AI:69:GLY:O	2.18	0.42
1:AA:1273:C:H2'	1:AA:1274:A:O4'	2.19	0.42
1:AA:234:C:H2'	1:AA:235:C:H6	1.84	0.42
1:AA:664:G:H5''	17:AR:52:ARG:NH2	2.34	0.42
1:AA:844:G:H2'	1:AA:845:A:H5''	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:AB:112:ARG:HH11	20:AB:112:ARG:HG3	1.83	0.42
3:AD:116:LEU:O	3:AD:121:ALA:HB3	2.19	0.42
4:AE:25:LYS:HE2	4:AE:25:LYS:C	2.39	0.42
5:AF:43:GLY:HA2	5:AF:58:HIS:CE1	2.53	0.42
6:AG:102:TRP:CD1	6:AG:102:TRP:N	2.86	0.42
7:AH:38:VAL:O	7:AH:42:GLU:HG2	2.19	0.42
8:AI:118:ARG:NH1	8:AI:122:ARG:HE	2.17	0.42
9:AJ:38:GLY:O	9:AJ:74:VAL:HA	2.19	0.42
13:AN:62:ARG:HA	13:AN:69:PRO:HA	2.01	0.42
1:AA:254:G:OP1	16:AQ:68:LYS:O	2.37	0.42
21:AU:26:GLY:C	21:AU:28:LEU:N	2.72	0.42
47:B0:11:LYS:HG3	47:B0:12:ARG:N	2.35	0.42
47:B0:7:PRO:O	47:B0:8:THR:C	2.58	0.42
50:B3:9:ALA:O	50:B3:13:PHE:HB3	2.19	0.42
23:BB:1017:G:O2'	23:BB:1018:U:H5'	2.20	0.42
23:BB:1064:C:H2'	23:BB:1065:U:O4'	2.19	0.42
23:BB:1236:G:O2'	23:BB:1237:A:P	2.77	0.42
23:BB:1270:C:H5''	23:BB:1271:G:O5'	2.19	0.42
23:BB:1370:C:H2'	23:BB:1371:G:O4'	2.18	0.42
23:BB:1408:G:H2'	23:BB:1409:U:C6	2.54	0.42
23:BB:1421:G:C2	23:BB:1422:G:C8	3.07	0.42
23:BB:1803:A:H2	23:BB:1823:G:O4'	2.02	0.42
23:BB:2361:G:O2'	23:BB:2362:C:H5'	2.19	0.42
23:BB:2643:G:H2'	23:BB:2644:G:O4'	2.18	0.42
23:BB:2649:C:H2'	23:BB:2650:U:C6	2.52	0.42
23:BB:647:G:H2'	23:BB:648:G:C8	2.55	0.42
23:BB:751:A:H5'	40:BS:90:LYS:HZ1	1.84	0.42
23:BB:795:C:H2'	23:BB:796:C:H6	1.85	0.42
23:BB:935:C:H2'	23:BB:936:A:C8	2.54	0.42
25:BC:171:VAL:HB	25:BC:182:LYS:CB	2.47	0.42
25:BC:124:LYS:O	25:BC:191:LEU:CD1	2.67	0.42
25:BC:68:ARG:CZ	25:BC:103:ILE:HD12	2.49	0.42
26:BD:37:VAL:HG22	26:BD:46:ARG:HD3	2.00	0.42
26:BD:61:THR:C	26:BD:63:PRO:HD2	2.40	0.42
27:BE:120:VAL:H	27:BE:189:THR:HA	1.84	0.42
27:BE:148:ILE:C	27:BE:148:ILE:HD12	2.38	0.42
27:BE:148:ILE:N	27:BE:183:PHE:HD2	2.17	0.42
27:BE:4:VAL:N	27:BE:117:ARG:HH21	2.17	0.42
28:BF:133:GLU:CG	28:BF:147:ARG:HG2	2.32	0.42
29:BG:150:TYR:O	29:BG:151:ARG:HB2	2.19	0.42
23:BB:1113:U:OP1	29:BG:1:SER:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:BH:114:GLU:HB3	30:BH:133:GLN:C	2.40	0.42
30:BH:4:ILE:N	30:BH:37:VAL:HB	2.19	0.42
30:BH:5:LEU:O	30:BH:6:LEU:HB3	2.19	0.42
31:BJ:100:VAL:CG2	31:BJ:101:ILE:H	2.26	0.42
31:BJ:55:ILE:HD13	31:BJ:132:HIS:NE2	2.35	0.42
31:BJ:72:LYS:HG3	31:BJ:90:GLU:HB2	2.01	0.42
32:BK:104:THR:OG1	32:BK:107:LEU:HB2	2.19	0.42
34:BM:33:LEU:HB2	34:BM:101:VAL:CG1	2.50	0.42
34:BM:35:ALA:C	34:BM:36:VAL:HG13	2.40	0.42
34:BM:90:GLU:O	34:BM:91:TYR:HB2	2.20	0.42
35:BN:28:LEU:CD2	35:BN:48:VAL:HG21	2.49	0.42
35:BN:51:LEU:HA	35:BN:54:LEU:HD13	2.00	0.42
37:BP:61:ARG:O	37:BP:62:LYS:C	2.57	0.42
37:BP:91:VAL:HB	37:BP:113:LEU:CG	2.47	0.42
39:BR:33:VAL:O	39:BR:33:VAL:HG12	2.19	0.42
41:BT:26:LYS:HD2	41:BT:26:LYS:C	2.40	0.42
42:BU:38:ILE:O	42:BU:62:ALA:HB3	2.19	0.42
24:BV:53:LYS:HE3	24:BV:55:GLU:OE2	2.19	0.42
43:BW:11:ASN:ND2	43:BW:13:ARG:NH2	2.67	0.42
44:BX:41:HIS:CD2	44:BX:41:HIS:H	2.36	0.42
1:CA:138:G:O2'	1:CA:139:A:H5'	2.19	0.42
1:CA:195:A:H1'	1:CA:222:C:O2'	2.18	0.42
1:CA:233:C:O2'	1:CA:234:C:H5'	2.19	0.42
1:CA:329:A:H8	1:CA:329:A:H5'	1.84	0.42
1:CA:626:G:H2'	1:CA:627:G:H8	1.83	0.42
1:CA:674:G:OP1	5:CF:51:ILE:HG13	2.20	0.42
1:CA:806:C:O2'	1:CA:807:A:H5'	2.20	0.42
1:CA:913:A:H4'	1:CA:914:A:H4'	2.01	0.42
20:CB:86:CYS:C	20:CB:88:GLN:N	2.73	0.42
2:CC:128:MET:O	2:CC:130:ARG:N	2.41	0.42
2:CC:139:ASN:O	2:CC:143:LEU:HD13	2.18	0.42
3:CD:24:VAL:HG23	3:CD:25:ARG:H	1.83	0.42
4:CE:102:THR:HG21	4:CE:123:LEU:HD21	2.01	0.42
7:CH:14:ARG:HG3	7:CH:15:ASN:N	2.33	0.42
9:CJ:52:LEU:CD1	9:CJ:62:ARG:HE	2.31	0.42
11:CL:100:ALA:O	11:CL:103:CYS:SG	2.78	0.42
13:CN:63:CYS:O	13:CN:67:GLY:HA2	2.18	0.42
15:CP:18:GLN:HA	15:CP:38:PHE:HA	2.02	0.42
19:CT:66:ILE:HG22	19:CT:67:HIS:N	2.34	0.42
48:D1:8:ILE:HG21	48:D1:27:ARG:HD3	2.01	0.42
48:D1:5:ARG:HA	48:D1:28:THR:HG1	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:DA:15:A:H1'	22:DA:109:A:C5	2.55	0.42
22:DA:56:G:H4'	22:DA:57:A:H5'	2.02	0.42
23:DB:1027:A:C2	23:DB:2488:G:H5''	2.53	0.42
23:DB:1213:A:H62	23:DB:1236:G:H1'	1.83	0.42
23:DB:125:A:C4'	49:D2:13:ASN:ND2	2.82	0.42
23:DB:1526:C:O2'	23:DB:1527:G:H5'	2.19	0.42
23:DB:2019:A:C6	23:DB:2020:A:N7	2.88	0.42
23:DB:2479:U:H6	23:DB:2479:U:O5'	2.02	0.42
23:DB:249:C:O3'	50:D3:7:ARG:HD3	2.19	0.42
23:DB:2633:G:H2'	23:DB:2634:A:O4'	2.19	0.42
23:DB:2671:G:O2'	23:DB:2672:U:H5'	2.19	0.42
23:DB:2686:G:H2'	23:DB:2687:U:O4'	2.19	0.42
23:DB:2705:A:H2'	23:DB:2706:A:O4'	2.19	0.42
23:DB:538:A:H2'	23:DB:539:G:O4'	2.19	0.42
23:DB:545:U:C4	23:DB:548:G:OP1	2.73	0.42
23:DB:678:C:H2'	23:DB:679:C:C6	2.53	0.42
23:DB:5:A:H2'	23:DB:6:A:H8	1.84	0.42
23:DB:906:U:H1'	34:DM:66:ARG:NH2	2.34	0.42
23:DB:1813:G:O4'	25:DC:45:ASN:HB3	2.20	0.42
25:DC:47:ARG:HB3	25:DC:48:ILE:H	1.54	0.42
27:DE:115:GLN:CB	27:DE:117:ARG:NH1	2.81	0.42
22:DA:43:C:C4'	28:DF:62:GLN:HE21	2.32	0.42
30:DH:108:VAL:CG1	30:DH:110:VAL:HB	2.49	0.42
30:DH:88:GLY:HA3	30:DH:125:THR:OG1	2.18	0.42
31:DJ:78:THR:H	31:DJ:84:ILE:HD13	1.84	0.42
23:DB:2561:U:O3'	32:DK:40:LYS:HE2	2.19	0.42
33:DL:71:ALA:O	33:DL:72:ALA:O	2.37	0.42
37:DP:18:SER:HB2	37:DP:87:ARG:NH1	2.34	0.42
38:DQ:52:ARG:O	38:DQ:53:LYS:C	2.58	0.42
41:DT:69:ARG:NH1	41:DT:69:ARG:HG2	2.34	0.42
42:DU:32:LYS:HA	42:DU:65:GLN:NE2	2.34	0.42
1:AA:1073:U:H2'	1:AA:1074:G:H8	1.84	0.42
1:AA:1073:U:C2	1:AA:1074:G:C8	3.08	0.42
1:AA:1178:G:N2	1:AA:1181:G:C8	2.87	0.42
1:AA:154:U:O2'	1:AA:155:A:H5'	2.20	0.42
1:AA:492:C:C3'	1:AA:493:A:H5''	2.49	0.42
4:AE:109:ALA:CB	4:AE:135:VAL:HG23	2.47	0.42
5:AF:29:ILE:HG21	5:AF:64:VAL:CG1	2.50	0.42
6:AG:37:THR:O	6:AG:41:ILE:HB	2.19	0.42
6:AG:90:VAL:HG23	6:AG:90:VAL:O	2.19	0.42
10:AK:62:ALA:O	10:AK:65:ALA:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AO:38:LEU:HB3	14:AO:42:PHE:CE1	2.55	0.42
47:B0:2:VAL:CG2	47:B0:3:GLN:H	2.33	0.42
50:B3:16:THR:CG2	50:B3:22:LYS:HG2	2.50	0.42
23:BB:651:G:OP1	50:B3:16:THR:HB	2.20	0.42
23:BB:1011:G:P	38:BQ:76:SER:HG	2.42	0.42
23:BB:1114:C:H2'	23:BB:1115:G:C8	2.55	0.42
23:BB:1221:C:O2'	23:BB:1222:U:H5'	2.19	0.42
23:BB:1401:G:H2'	23:BB:1402:U:C6	2.54	0.42
23:BB:1453:A:H62	35:BN:72:ASP:CG	2.23	0.42
23:BB:1498:C:H2'	23:BB:1499:C:C6	2.54	0.42
23:BB:1513:U:O2'	23:BB:1514:G:H5'	2.19	0.42
23:BB:1674:G:N2	23:BB:1677:A:N1	2.68	0.42
23:BB:1685:C:H2'	23:BB:1686:C:C6	2.54	0.42
23:BB:1791:A:H1'	25:BC:206:LYS:HZ1	1.81	0.42
23:BB:1804:C:P	25:BC:254:LYS:HG2	2.58	0.42
23:BB:1891:G:H2'	23:BB:1892:C:H6	1.84	0.42
23:BB:2559:C:O2'	23:BB:2560:A:H5'	2.19	0.42
23:BB:2705:A:H2'	23:BB:2706:A:O4'	2.19	0.42
23:BB:2811:G:H2'	23:BB:2812:G:C8	2.54	0.42
23:BB:2813:A:O2'	23:BB:2814:A:H5'	2.18	0.42
23:BB:312:G:C2	23:BB:313:G:C8	3.07	0.42
23:BB:748:G:OP2	40:BS:88:ARG:HD3	2.19	0.42
23:BB:970:U:H2'	23:BB:971:G:C8	2.55	0.42
25:BC:170:TYR:HA	25:BC:185:ALA:HB2	2.01	0.42
26:BD:205:PRO:O	26:BD:206:ALA:HB2	2.19	0.42
27:BE:53:THR:O	27:BE:55:SER:N	2.52	0.42
30:BH:124:THR:HG22	30:BH:125:THR:N	2.35	0.42
52:BI:63:ASP:O	52:BI:65:SER:N	2.53	0.42
31:BJ:68:LYS:CA	31:BJ:71:ASP:OD1	2.65	0.42
1:AA:1422:G:P	32:BK:54:LYS:HZ1	2.43	0.42
33:BL:58:TYR:N	33:BL:58:TYR:CD2	2.84	0.42
34:BM:3:GLN:CG	34:BM:46:ILE:HB	2.43	0.42
34:BM:86:LYS:HB3	34:BM:87:GLY:H	1.67	0.42
35:BN:28:LEU:CD2	35:BN:113:ILE:HG21	2.49	0.42
37:BP:47:ILE:HG22	37:BP:48:ALA:N	2.17	0.42
37:BP:77:SER:N	37:BP:78:PRO:CD	2.73	0.42
39:BR:101:ILE:HB	39:BR:102:SER:H	1.69	0.42
23:BB:1225:G:C5'	39:BR:78:ARG:HH22	2.27	0.42
41:BT:64:LYS:HZ2	41:BT:79:ASP:HB2	1.85	0.42
42:BU:21:ARG:NH1	42:BU:21:ARG:HG2	2.33	0.42
23:BB:500:G:OP1	42:BU:40:LEU:HG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:BW:35:ILE:HG13	43:BW:36:ILE:N	2.33	0.42
43:BW:44:PHE:HB3	43:BW:76:ARG:CB	2.49	0.42
1:CA:1141:C:H2'	1:CA:1142:G:H8	1.83	0.42
1:CA:1306:A:C2'	1:CA:1307:U:H5'	2.49	0.42
1:CA:1394:A:H8	1:CA:1394:A:OP1	2.03	0.42
1:CA:349:A:H2'	1:CA:350:G:C8	2.54	0.42
1:CA:538:G:H2'	1:CA:539:A:H8	1.85	0.42
1:CA:838:G:H2'	1:CA:839:C:O4'	2.20	0.42
1:CA:976:G:OP2	1:CA:1358:U:O2'	2.32	0.42
20:CB:15:PHE:HB2	20:CB:16:GLY:H	1.65	0.42
20:CB:61:SER:OG	20:CB:62:ARG:HD2	2.19	0.42
20:CB:93:HIS:O	20:CB:94:ARG:C	2.57	0.42
2:CC:130:ARG:HD2	2:CC:130:ARG:HA	1.86	0.42
2:CC:172:VAL:O	2:CC:174:LEU:HD12	2.19	0.42
2:CC:183:TYR:HA	2:CC:199:VAL:O	2.19	0.42
2:CC:182:ASP:CB	2:CC:203:LYS:HE2	2.43	0.42
3:CD:159:GLU:O	3:CD:162:GLU:HG2	2.19	0.42
4:CE:105:ILE:HD12	4:CE:123:LEU:CD2	2.50	0.42
5:CF:74:LEU:HD13	5:CF:74:LEU:C	2.39	0.42
8:CI:49:GLN:O	8:CI:53:LEU:HD23	2.19	0.42
19:CT:79:THR:O	19:CT:82:ILE:HG12	2.19	0.42
47:D0:27:LEU:O	47:D0:38:LEU:HD13	2.19	0.42
33:DL:63:LYS:CB	50:D3:26:ALA:HB2	2.47	0.42
23:DB:2393:U:C5	50:D3:30:HIS:HE1	2.37	0.42
50:D3:57:VAL:HG22	50:D3:57:VAL:O	2.19	0.42
23:DB:1098:A:HO2'	52:DI:4:VAL:C	2.21	0.42
23:DB:1220:G:H2'	23:DB:1221:C:H6	1.83	0.42
23:DB:1252:G:H21	38:DQ:32:ARG:CZ	2.32	0.42
23:DB:1401:G:H2'	23:DB:1402:U:C6	2.54	0.42
23:DB:2082:A:N6	23:DB:2237:G:H1'	2.34	0.42
23:DB:2248:C:H3'	23:DB:2249:U:C6	2.55	0.42
23:DB:954:G:O2'	23:DB:2274:A:N1	2.44	0.42
23:DB:2408:U:H2'	23:DB:2409:G:C8	2.54	0.42
23:DB:2259:U:C6	23:DB:2427:C:C4	3.07	0.42
23:DB:2462:C:H2'	23:DB:2463:C:H6	1.84	0.42
23:DB:2570:G:O2'	23:DB:2571:U:H5'	2.20	0.42
23:DB:539:G:H2'	23:DB:540:C:H6	1.83	0.42
23:DB:662:G:H5'	33:DL:26:GLY:H	1.83	0.42
23:DB:756:A:H2'	23:DB:757:G:O4'	2.20	0.42
25:DC:163:ILE:HG12	25:DC:173:LEU:HD23	2.01	0.42
25:DC:245:THR:O	25:DC:246:PRO:C	2.57	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2821:A:OP2	26:DD:115:GLY:O	2.38	0.42
26:DD:48:ILE:HG22	26:DD:49:GLN:H	1.80	0.42
27:DE:112:LEU:O	27:DE:113:VAL:HB	2.20	0.42
28:DF:50:ASP:C	28:DF:52:ALA:H	2.22	0.42
29:DG:57:TYR:O	29:DG:58:ALA:C	2.56	0.42
30:DH:132:PHE:HE2	30:DH:134:VAL:HG22	1.84	0.42
30:DH:35:LYS:HE2	30:DH:35:LYS:HB2	1.91	0.42
30:DH:4:ILE:HD11	30:DH:37:VAL:CA	2.48	0.42
31:DJ:15:TRP:CB	31:DJ:139:VAL:HA	2.49	0.42
31:DJ:7:LYS:N	31:DJ:8:PRO:CD	2.82	0.42
32:DK:43:ILE:HG12	32:DK:52:VAL:CG1	2.46	0.42
33:DL:77:ILE:O	33:DL:110:VAL:O	2.37	0.42
33:DL:81:ASP:O	33:DL:85:VAL:HG12	2.19	0.42
34:DM:41:LEU:CD2	34:DM:46:ILE:HD11	2.49	0.42
36:DO:82:ALA:HB1	36:DO:87:ILE:HG21	2.01	0.42
37:DP:7:LEU:HA	37:DP:10:GLU:CD	2.39	0.42
23:DB:533:G:N3	38:DQ:40:LYS:CG	2.82	0.42
38:DQ:44:TYR:O	38:DQ:48:ASP:N	2.51	0.42
38:DQ:52:ARG:HD3	38:DQ:53:LYS:HZ1	1.83	0.42
38:DQ:57:ARG:NH2	38:DQ:92:LYS:HZ3	2.16	0.42
23:DB:2010:G:H5''	40:DS:42:LYS:HB3	2.00	0.42
24:DV:9:ARG:CZ	24:DV:20:LEU:HD11	2.50	0.42
43:DW:67:LYS:CD	43:DW:69:GLU:H	2.33	0.42
45:DY:43:ILE:O	45:DY:47:ILE:HG13	2.20	0.42
1:AA:1065:U:C1'	1:AA:1066:C:OP2	2.66	0.42
1:AA:123:U:H2'	1:AA:124:C:C6	2.54	0.42
1:AA:128:G:O2'	1:AA:129:A:H5'	2.18	0.42
1:AA:1306:A:C2	1:AA:1332:A:N9	2.87	0.42
1:AA:138:G:O2'	1:AA:139:A:H5'	2.19	0.42
1:AA:329:A:H5'	1:AA:329:A:H8	1.84	0.42
1:AA:343:U:O2'	1:AA:344:A:H2'	2.20	0.42
1:AA:442:G:H2'	1:AA:443:C:H6	1.84	0.42
1:AA:913:A:H4'	1:AA:914:A:O5'	2.19	0.42
3:AD:173:ASP:HB3	3:AD:178:GLU:HB3	2.01	0.42
3:AD:191:SER:O	3:AD:192:ALA:CB	2.68	0.42
1:AA:15:G:O2'	4:AE:28:ARG:HD2	2.20	0.42
6:AG:45:ALA:HB1	6:AG:120:ALA:HB2	2.00	0.42
13:AN:12:ARG:HH11	13:AN:60:ARG:HH12	1.64	0.42
18:AS:35:ARG:HG2	18:AS:50:VAL:CG1	2.49	0.42
1:AA:1314:C:H5	18:AS:5:LYS:HZ3	1.63	0.42
47:B0:26:SER:HB2	47:B0:38:LEU:CD1	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:B1:31:GLU:C	48:B1:32:LYS:HZ2	2.22	0.42
48:B1:51:ALA:HB3	48:B1:52:LYS:HD2	2.01	0.42
51:B4:12:ARG:CD	51:B4:12:ARG:H	2.06	0.42
23:BB:1173:U:H2'	23:BB:1174:U:H4'	2.00	0.42
23:BB:1266:G:O2'	23:BB:1267:U:OP2	2.38	0.42
23:BB:126:A:O5'	49:B2:19:ARG:HB3	2.19	0.42
23:BB:1298:C:O2'	23:BB:1301:A:O2'	2.36	0.42
23:BB:1604:C:H5''	55:BB:3376:HOH:O	2.18	0.42
23:BB:1770:G:O2'	23:BB:1771:C:H5'	2.19	0.42
23:BB:2263:C:H4'	23:BB:2329:U:H4'	2.01	0.42
23:BB:2463:C:O2'	23:BB:2464:G:H5'	2.19	0.42
23:BB:299:A:N6	23:BB:322:A:O2'	2.49	0.42
23:BB:453:A:H4'	23:BB:472:A:H62	1.84	0.42
23:BB:679:C:O2'	23:BB:680:C:H5'	2.19	0.42
23:BB:693:A:H2'	23:BB:694:U:C6	2.54	0.42
23:BB:919:U:H6	23:BB:919:U:O5'	2.03	0.42
23:BB:84:A:N6	23:BB:99:U:H5'	2.35	0.42
25:BC:207:ALA:O	25:BC:208:GLY:C	2.57	0.42
23:BB:1566:A:C6	25:BC:213:ARG:CZ	3.02	0.42
27:BE:112:LEU:O	27:BE:116:ASP:N	2.52	0.42
27:BE:156:ASN:CG	27:BE:157:LEU:H	2.23	0.42
27:BE:68:ALA:O	27:BE:69:ARG:HB3	2.20	0.42
27:BE:74:LYS:O	27:BE:75:SER:C	2.58	0.42
29:BG:99:GLY:C	29:BG:101:VAL:H	2.22	0.42
30:BH:90:LEU:HD13	30:BH:123:ARG:O	2.19	0.42
30:BH:89:LYS:HZ1	30:BH:90:LEU:HB2	1.85	0.42
52:BI:21:PRO:HB2	52:BI:22:PRO:CD	2.44	0.42
52:BI:92:PRO:C	52:BI:94:LYS:H	2.23	0.42
31:BJ:124:VAL:O	31:BJ:125:TYR:HB2	2.18	0.42
31:BJ:12:LYS:O	31:BJ:13:ARG:HG3	2.19	0.42
32:BK:2:ILE:N	32:BK:33:ALA:N	2.68	0.42
32:BK:71:ARG:HH11	32:BK:71:ARG:HG2	1.84	0.42
33:BL:17:LYS:HD3	33:BL:17:LYS:HA	1.92	0.42
33:BL:78:ARG:O	33:BL:78:ARG:NH1	2.53	0.42
34:BM:3:GLN:HE21	34:BM:3:GLN:HB2	1.67	0.42
37:BP:61:ARG:CB	37:BP:61:ARG:NH1	2.81	0.42
31:BJ:45:THR:CB	38:BQ:63:ARG:HH22	2.16	0.42
38:BQ:63:ARG:HG3	38:BQ:64:ILE:N	2.33	0.42
38:BQ:89:ILE:O	38:BQ:89:ILE:HG23	2.20	0.42
39:BR:41:ILE:HD13	39:BR:43:ASN:HA	2.00	0.42
39:BR:51:VAL:HG13	39:BR:52:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BU:70:ALA:O	42:BU:72:PHE:N	2.53	0.42
43:BW:36:ILE:HD12	43:BW:69:GLU:CD	2.39	0.42
45:BY:22:THR:O	45:BY:26:LEU:N	2.53	0.42
46:BZ:2:LYS:HB2	46:BZ:3:LYS:H	1.69	0.42
1:CA:113:G:O4'	1:CA:354:G:H4'	2.19	0.42
1:CA:1164:G:H2'	1:CA:1165:U:H6	1.85	0.42
1:CA:1237:C:O4'	1:CA:1334:G:N2	2.50	0.42
1:CA:1250:A:H5''	8:CI:68:GLY:HA2	2.01	0.42
1:CA:1252:A:H4'	1:CA:1369:C:H4'	2.01	0.42
1:CA:167:A:O2'	1:CA:168:G:H5'	2.19	0.42
1:CA:6:G:H4'	1:CA:298:A:H4'	2.01	0.42
1:CA:492:C:C3'	1:CA:493:A:H5''	2.50	0.42
1:CA:64:G:H4'	1:CA:65:A:H5''	2.01	0.42
20:CB:86:CYS:HB3	20:CB:88:GLN:CG	2.49	0.42
4:CE:10:LEU:HD12	4:CE:10:LEU:O	2.20	0.42
5:CF:15:SER:HA	5:CF:18:VAL:HG23	2.01	0.42
5:CF:84:VAL:HG22	5:CF:85:ILE:H	1.85	0.42
6:CG:49:LEU:C	6:CG:51:GLN:N	2.73	0.42
8:CI:19:PHE:O	8:CI:62:LEU:HA	2.19	0.42
8:CI:94:ARG:HG3	8:CI:98:ARG:NH2	2.33	0.42
9:CJ:66:GLU:HG2	9:CJ:67:ILE:N	2.34	0.42
11:CL:66:ILE:HG21	11:CL:71:HIS:HB3	2.00	0.42
12:CM:21:ILE:O	12:CM:22:TYR:C	2.57	0.42
21:CU:36:PHE:HB2	21:CU:39:LYS:CB	2.47	0.42
49:D2:21:ARG:HG3	49:D2:31:LEU:CD1	2.46	0.42
49:D2:33:ARG:O	49:D2:37:LYS:HB3	2.20	0.42
23:DB:1189:A:H2'	23:DB:1190:G:H5'	2.02	0.42
23:DB:1266:G:O2'	23:DB:1267:U:OP2	2.38	0.42
23:DB:1290:C:O2'	23:DB:1291:C:H5'	2.19	0.42
23:DB:1294:U:C4	23:DB:1295:C:C5	3.07	0.42
23:DB:1314:C:H2'	23:DB:1314:C:O2	2.18	0.42
23:DB:1592:C:H2'	23:DB:1593:A:H8	1.84	0.42
23:DB:1829:A:H3'	23:DB:1830:C:C6	2.55	0.42
23:DB:2153:C:H2'	23:DB:2154:A:H8	1.84	0.42
23:DB:2259:U:O2'	23:DB:2260:C:H5'	2.19	0.42
23:DB:2282:G:O2'	23:DB:2283:C:OP2	2.37	0.42
23:DB:222:A:N1	23:DB:233:A:H5''	2.35	0.42
23:DB:2723:C:H2'	23:DB:2724:U:O4'	2.20	0.42
23:DB:62:U:H3'	23:DB:63:A:C8	2.54	0.42
25:DC:68:ARG:NE	25:DC:128:THR:HG23	2.34	0.42
25:DC:107:LYS:CG	25:DC:194:VAL:HG11	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:241:LYS:HB3	25:DC:243:PRO:HD3	2.02	0.42
25:DC:93:VAL:HG21	25:DC:103:ILE:HG13	2.00	0.42
26:DD:168:GLU:H	26:DD:168:GLU:CD	2.23	0.42
27:DE:126:VAL:HG21	27:DE:132:LYS:HZ2	1.82	0.42
28:DF:92:GLY:O	28:DF:95:MET:HG2	2.19	0.42
52:DI:129:GLU:O	52:DI:133:ARG:HG3	2.19	0.42
31:DJ:135:GLN:HE21	31:DJ:138:GLN:N	2.14	0.42
31:DJ:35:ARG:HA	31:DJ:35:ARG:NE	2.35	0.42
31:DJ:61:LYS:HE2	31:DJ:69:ARG:HH22	1.84	0.42
31:DJ:84:ILE:CD1	31:DJ:85:LYS:N	2.73	0.42
35:DN:37:THR:CG2	35:DN:39:PRO:HD2	2.41	0.42
36:DO:38:GLN:O	36:DO:40:ILE:HD13	2.18	0.42
36:DO:25:ARG:HH21	36:DO:94:ARG:NH1	2.17	0.42
37:DP:105:LYS:HE2	37:DP:105:LYS:HB3	1.81	0.42
41:DT:11:LEU:HB2	41:DT:12:ARG:H	1.59	0.42
41:DT:69:ARG:HH11	41:DT:69:ARG:HG2	1.85	0.42
42:DU:29:SER:O	42:DU:30:SER:HB2	2.19	0.42
23:DB:372:G:N7	46:DZ:57:VAL:HG21	2.34	0.42
1:AA:1299:A:N6	1:AA:1302:C:H5	2.15	0.42
1:AA:1490:U:H3'	1:AA:1491:G:H8	1.85	0.42
1:AA:302:G:H2'	1:AA:303:A:H8	1.83	0.42
1:AA:41:G:H2'	1:AA:42:G:C8	2.54	0.42
1:AA:838:G:H2'	1:AA:839:C:O4'	2.20	0.42
1:AA:868:C:H2'	1:AA:869:G:O4'	2.20	0.42
20:AB:133:ALA:O	20:AB:137:THR:HG23	2.20	0.42
2:AC:120:THR:HA	2:AC:123:LEU:HD12	2.01	0.42
2:AC:54:ILE:HG23	2:AC:54:ILE:O	2.19	0.42
8:AI:24:ASN:HA	8:AI:26:LYS:NZ	2.34	0.42
9:AJ:48:ARG:HB3	9:AJ:66:GLU:HB3	2.02	0.42
10:AK:52:ARG:O	10:AK:54:SER:N	2.51	0.42
16:AQ:17:GLU:O	16:AQ:18:LYS:HB2	2.18	0.42
16:AQ:56:ASP:CG	16:AQ:80:LYS:HA	2.39	0.42
47:B0:8:THR:HG23	47:B0:9:ARG:N	2.26	0.42
22:BA:54:G:H2'	22:BA:55:U:H6	1.83	0.42
23:BB:1047:G:HO2'	23:BB:1048:A:P	2.42	0.42
23:BB:1250:G:C5'	38:BQ:5:ARG:HG2	2.49	0.42
23:BB:2223:G:C2'	23:BB:2224:G:H5'	2.50	0.42
23:BB:2286:G:H3'	23:BB:2286:G:N3	2.34	0.42
23:BB:2787:C:O2'	23:BB:2788:C:H5'	2.19	0.42
23:BB:2840:C:C2	23:BB:2841:C:C5	3.07	0.42
23:BB:63:A:H8	23:BB:63:A:OP2	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:649:G:H2'	23:BB:650:C:C6	2.55	0.42
23:BB:64:A:H2'	23:BB:65:U:H6	1.80	0.42
23:BB:723:C:H2'	23:BB:724:U:C6	2.55	0.42
25:BC:17:LYS:N	25:BC:17:LYS:HZ2	2.17	0.42
25:BC:31:PRO:O	25:BC:32:LEU:O	2.37	0.42
26:BD:1:MET:HB3	26:BD:84:LEU:HB2	2.00	0.42
27:BE:195:GLN:H	27:BE:199:MET:CA	2.28	0.42
29:BG:104:LEU:HB3	29:BG:106:LEU:CD1	2.49	0.42
29:BG:51:PHE:CD2	29:BG:64:ALA:HA	2.54	0.42
30:BH:2:GLN:HE22	30:BH:19:VAL:C	2.23	0.42
31:BJ:93:ILE:HD13	31:BJ:100:VAL:HG11	2.01	0.42
31:BJ:3:THR:HG21	38:BQ:93:ILE:CG1	2.40	0.42
32:BK:38:ILE:HD13	32:BK:61:VAL:HG12	2.00	0.42
32:BK:76:VAL:HG12	32:BK:77:ILE:N	2.35	0.42
34:BM:11:LYS:HD3	34:BM:11:LYS:N	2.34	0.42
36:BO:41:ALA:CB	36:BO:45:SER:HB3	2.49	0.42
36:BO:59:ALA:CB	36:BO:63:LYS:HE2	2.50	0.42
32:BK:80:ASP:OD1	37:BP:71:ARG:HG2	2.19	0.42
41:BT:15:HIS:HB2	41:BT:31:VAL:CG2	2.50	0.42
42:BU:54:PRO:CG	42:BU:55:GLY:H	2.30	0.42
24:BV:6:ALA:HB1	24:BV:41:GLU:O	2.19	0.42
24:BV:2:PHE:HE2	24:BV:55:GLU:HB3	1.84	0.42
24:BV:80:HIS:HB2	24:BV:85:LYS:HG3	2.01	0.42
23:BB:2080:A:OP1	46:BZ:18:CYS:HB2	2.20	0.42
1:CA:1217:C:O2'	1:CA:1218:C:H5'	2.20	0.42
1:CA:738:C:H2'	1:CA:739:C:C6	2.55	0.42
1:CA:75:G:H2'	1:CA:76:G:O4'	2.19	0.42
1:CA:926:G:H5'	1:CA:927:G:C5'	2.49	0.42
1:CA:997:U:H2'	1:CA:998:C:H6	1.79	0.42
20:CB:99:MET:CA	20:CB:106:VAL:HG21	2.47	0.42
20:CB:163:ILE:HD11	20:CB:209:VAL:HG11	2.01	0.42
2:CC:128:MET:SD	2:CC:132:ALA:N	2.92	0.42
3:CD:33:ILE:HG12	3:CD:34:GLU:N	2.34	0.42
3:CD:10:LEU:HD22	3:CD:62:ARG:CZ	2.49	0.42
4:CE:40:ASP:CG	4:CE:44:ARG:HB2	2.39	0.42
4:CE:52:ALA:HB3	4:CE:58:ALA:HB2	2.01	0.42
6:CG:60:ALA:O	6:CG:64:ALA:HB3	2.19	0.42
7:CH:72:GLU:N	7:CH:129:ALA:HB2	2.27	0.42
9:CJ:18:ILE:HD13	9:CJ:72:ARG:HG2	2.00	0.42
14:CO:63:ARG:NH2	14:CO:87:ARG:NH1	2.67	0.42
15:CP:54:LEU:CD1	15:CP:80:LYS:HG2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:CS:73:PHE:CD1	18:CS:73:PHE:N	2.88	0.42
47:D0:11:LYS:HA	47:D0:11:LYS:HD3	1.90	0.42
47:D0:42:ILE:HG21	47:D0:45:ASP:OD2	2.19	0.42
49:D2:39:ARG:HD2	49:D2:39:ARG:N	2.35	0.42
23:DB:1039:A:H2'	23:DB:1040:A:C8	2.53	0.42
23:DB:1298:C:O2'	23:DB:1301:A:O2'	2.36	0.42
23:DB:1683:U:O2'	23:DB:1684:G:H5'	2.19	0.42
23:DB:1765:U:O2'	23:DB:1766:G:H5'	2.19	0.42
23:DB:2047:C:H2'	23:DB:2048:G:C8	2.52	0.42
23:DB:2267:A:H3'	23:DB:2267:A:N3	2.34	0.42
23:DB:2543:G:H2'	23:DB:2544:G:O4'	2.20	0.42
23:DB:2733:A:O5'	23:DB:2733:A:H8	2.03	0.42
23:DB:361:G:H2'	23:DB:362:A:C8	2.51	0.42
23:DB:448:U:H3'	27:DE:79:ARG:CZ	2.49	0.42
23:DB:975:A:H2'	23:DB:975:A:N3	2.35	0.42
25:DC:175:LEU:H	25:DC:175:LEU:HG	1.67	0.42
25:DC:253:GLY:H	25:DC:254:LYS:HE3	1.83	0.42
25:DC:265:PHE:O	25:DC:266:ILE:HG12	2.20	0.42
23:DB:2572:A:OP2	26:DD:149:ASN:HB3	2.19	0.42
22:DA:45:A:O4'	28:DF:91:ARG:CZ	2.67	0.42
29:DG:124:CYS:C	29:DG:126:THR:H	2.21	0.42
29:DG:124:CYS:HB3	29:DG:130:ILE:CD1	2.49	0.42
29:DG:171:LYS:HZ3	29:DG:174:LYS:HD2	1.85	0.42
29:DG:16:VAL:HG12	29:DG:17:LYS:H	1.84	0.42
30:DH:103:VAL:O	30:DH:106:ALA:HB3	2.20	0.42
30:DH:129:GLU:HG3	30:DH:129:GLU:H	1.57	0.42
30:DH:54:LEU:HD23	30:DH:57:LYS:CD	2.48	0.42
52:DI:37:PHE:HB2	52:DI:66:PHE:CE2	2.55	0.42
31:DJ:69:ARG:NH1	31:DJ:69:ARG:HG3	2.31	0.42
32:DK:15:GLY:CA	32:DK:52:VAL:HB	2.49	0.42
23:DB:1667:G:OP1	32:DK:6:THR:HA	2.20	0.42
33:DL:124:GLY:H	33:DL:142:ILE:HG13	1.85	0.42
36:DO:67:ASN:HD22	36:DO:68:LYS:H	1.68	0.42
36:DO:94:ARG:HG2	36:DO:94:ARG:HH11	1.84	0.42
32:DK:78:ARG:CG	37:DP:72:VAL:HG21	2.18	0.42
38:DQ:17:LEU:HD12	38:DQ:17:LEU:HA	1.89	0.42
38:DQ:50:ARG:NH1	38:DQ:53:LYS:HE3	2.34	0.42
38:DQ:94:LEU:C	38:DQ:94:LEU:HD12	2.39	0.42
38:DQ:94:LEU:O	38:DQ:98:ALA:N	2.49	0.42
39:DR:65:ALA:N	39:DR:100:GLY:HA2	2.34	0.42
41:DT:6:ARG:O	41:DT:8:LEU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DU:42:LYS:H	42:DU:57:ILE:CG2	2.32	0.42
46:DZ:39:LYS:HD3	46:DZ:61:ASN:CG	2.39	0.42
1:AA:1432:G:O2'	1:AA:1468:A:N6	2.50	0.42
1:AA:1447:A:H5'	1:AA:1448:C:H5	1.85	0.42
1:AA:1521:C:H2'	1:AA:1522:U:H6	1.84	0.42
1:AA:278:G:H21	1:AA:279:A:N6	2.10	0.42
1:AA:512:U:O2'	1:AA:513:C:H5'	2.19	0.42
1:AA:687:A:C2	1:AA:704:A:C6	3.07	0.42
1:AA:908:A:O2'	1:AA:909:A:H5'	2.19	0.42
20:AB:15:PHE:HD1	20:AB:16:GLY:N	2.18	0.42
6:AG:20:GLU:O	6:AG:23:ALA:HB3	2.19	0.42
7:AH:83:ARG:NH1	7:AH:123:GLU:OE1	2.53	0.42
9:AJ:25:ILE:HG23	9:AJ:29:ALA:CB	2.47	0.42
10:AK:70:ALA:HA	10:AK:74:LYS:HD3	2.01	0.42
11:AL:43:LYS:N	11:AL:44:PRO:CD	2.81	0.42
1:AA:529:G:O6	11:AL:45:ASN:HA	2.19	0.42
12:AM:10:ASP:HA	12:AM:44:ILE:HD11	2.01	0.42
12:AM:23:GLY:CA	12:AM:64:VAL:HG13	2.43	0.42
16:AQ:30:HIS:N	16:AQ:35:LYS:H	2.17	0.42
16:AQ:7:LEU:HB2	16:AQ:60:ILE:HG12	2.02	0.42
17:AR:19:GLU:C	17:AR:21:ASP:H	2.22	0.42
17:AR:23:LYS:C	17:AR:25:ILE:H	2.22	0.42
1:AA:958:A:C4'	18:AS:54:ARG:HH21	2.33	0.42
47:B0:36:LYS:HA	47:B0:41:HIS:C	2.40	0.42
23:BB:1068:G:C6	23:BB:1069:A:N6	2.87	0.42
23:BB:1403:A:H2'	23:BB:1404:C:C6	2.54	0.42
23:BB:1713:A:N6	23:BB:1745:A:H61	2.18	0.42
23:BB:1637:A:H5'	23:BB:1760:C:O2'	2.20	0.42
23:BB:181:A:O2'	23:BB:182:A:H5'	2.19	0.42
23:BB:1841:U:C2	23:BB:1842:G:C8	3.07	0.42
23:BB:2347:C:H4'	23:BB:2347:C:OP1	2.19	0.42
23:BB:2540:C:O2'	23:BB:2541:A:H5'	2.20	0.42
23:BB:2686:G:H2'	23:BB:2687:U:C6	2.54	0.42
23:BB:2723:C:H2'	23:BB:2724:U:O4'	2.20	0.42
23:BB:26:G:C2'	23:BB:27:G:O4'	2.67	0.42
23:BB:678:C:H2'	23:BB:679:C:C6	2.53	0.42
23:BB:775:G:C5	23:BB:794:A:C8	3.08	0.42
26:BD:175:LEU:CD2	26:BD:190:LYS:HB3	2.50	0.42
31:BJ:132:HIS:C	31:BJ:134:ALA:H	2.22	0.42
33:BL:34:GLY:O	33:BL:35:HIS:C	2.57	0.42
33:BL:78:ARG:NH1	33:BL:80:SER:N	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BM:69:PRO:HA	34:BM:93:VAL:HB	2.01	0.42
35:BN:12:ARG:HD3	35:BN:20:MET:CE	2.50	0.42
35:BN:13:ASN:C	35:BN:15:SER:H	2.23	0.42
23:BB:1156:A:C5	38:BQ:47:ARG:NH1	2.88	0.42
39:BR:21:ARG:HH11	39:BR:21:ARG:HG2	1.84	0.42
40:BS:50:VAL:HG21	40:BS:103:ILE:CG2	2.50	0.42
40:BS:29:VAL:HG21	40:BS:70:LYS:N	2.35	0.42
40:BS:82:MET:HB2	40:BS:98:LYS:CG	2.50	0.42
42:BU:33:VAL:HG12	42:BU:35:VAL:HG23	2.02	0.42
42:BU:90:LYS:HD2	42:BU:91:LYS:N	2.34	0.42
46:BZ:15:SER:HB3	46:BZ:23:LYS:HE2	2.02	0.42
1:CA:1014:A:H2'	1:CA:1015:G:C8	2.54	0.42
1:CA:1082:A:H2'	1:CA:1083:U:O4'	2.20	0.42
1:CA:1200:C:OP1	1:CA:1201:A:H3'	2.20	0.42
1:CA:109:A:N1	1:CA:326:G:C6	2.88	0.42
1:CA:404:G:H2'	1:CA:405:U:H6	1.85	0.42
1:CA:41:G:H2'	1:CA:42:G:C8	2.54	0.42
1:CA:429:U:H4'	1:CA:430:A:O5'	2.16	0.42
1:CA:538:G:H2'	1:CA:539:A:C8	2.54	0.42
1:CA:69:G:H2'	1:CA:70:U:C6	2.55	0.42
1:CA:896:C:O2'	1:CA:897:C:H5'	2.19	0.42
2:CC:110:LEU:HD13	2:CC:143:LEU:CD2	2.50	0.42
3:CD:195:ASN:HB2	3:CD:198:LEU:HD12	2.02	0.42
5:CF:29:ILE:HG21	5:CF:64:VAL:CG1	2.49	0.42
7:CH:65:PHE:CD2	7:CH:66:GLN:HG2	2.55	0.42
12:CM:70:ARG:NH1	28:DF:135:ILE:HD13	2.35	0.42
12:CM:78:ARG:HD2	12:CM:79:LEU:N	2.35	0.42
14:CO:87:ARG:O	14:CO:87:ARG:HG3	2.20	0.42
15:CP:10:GLY:HA3	15:CP:15:PRO:HA	2.01	0.42
15:CP:26:ASN:ND2	15:CP:31:ARG:HB3	2.34	0.42
23:DB:1264:A:H5'	47:D0:7:PRO:HG3	2.01	0.42
23:DB:686:U:O2'	49:D2:5:PHE:O	2.34	0.42
50:D3:44:ARG:HG2	50:D3:45:PRO:HD2	2.02	0.42
23:DB:249:C:O2'	50:D3:7:ARG:NE	2.52	0.42
51:D4:24:ARG:HE	51:D4:37:GLN:HA	1.84	0.42
23:DB:1029:A:H3'	23:DB:1030:C:C6	2.53	0.42
23:DB:1372:U:H2'	23:DB:1373:A:C8	2.55	0.42
23:DB:164:C:H2'	23:DB:165:A:H5'	2.01	0.42
23:DB:1773:A:H2'	23:DB:1774:C:O4'	2.19	0.42
23:DB:2135:A:N3	23:DB:2135:A:H2'	2.35	0.42
23:DB:2282:G:H5''	23:DB:2283:C:O4'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2425:A:H4'	23:DB:2426:A:OP2	2.19	0.42
23:DB:2540:C:O2'	23:DB:2541:A:H5'	2.20	0.42
23:DB:506:G:H1'	23:DB:507:A:C8	2.54	0.42
23:DB:647:G:H2'	23:DB:648:G:C8	2.55	0.42
23:DB:795:C:H2'	23:DB:796:C:C6	2.55	0.42
23:DB:982:C:O2	23:DB:982:C:H2'	2.20	0.42
25:DC:110:LYS:HE3	25:DC:110:LYS:CA	2.49	0.42
25:DC:21:PRO:N	25:DC:202:ARG:HH11	2.18	0.42
25:DC:222:THR:OG1	25:DC:223:ALA:N	2.53	0.42
25:DC:42:ARG:HE	25:DC:43:ASN:H	1.68	0.42
25:DC:8:THR:O	25:DC:10:PRO:HD3	2.20	0.42
26:DD:7:LYS:HA	26:DD:26:VAL:HA	2.01	0.42
26:DD:46:ARG:N	26:DD:82:PHE:HA	2.35	0.42
27:DE:175:ILE:CG2	27:DE:180:LEU:HD11	2.49	0.42
28:DF:111:ARG:HG2	28:DF:111:ARG:HH11	1.85	0.42
28:DF:65:LEU:O	28:DF:66:ILE:HB	2.20	0.42
29:DG:36:LEU:HB2	29:DG:40:VAL:HG11	2.02	0.42
29:DG:43:LYS:HB3	29:DG:44:HIS:H	1.71	0.42
29:DG:4:ALA:HA	29:DG:68:ARG:HH22	1.85	0.42
52:DI:10:LEU:C	52:DI:10:LEU:HD12	2.39	0.42
52:DI:121:ILE:HD11	52:DI:122:GLU:OE2	2.20	0.42
31:DJ:74:TYR:OH	31:DJ:100:VAL:HG23	2.20	0.42
23:DB:906:U:H4'	34:DM:26:VAL:CG1	2.49	0.42
35:DN:67:PHE:O	35:DN:71:ARG:HA	2.20	0.42
37:DP:100:ARG:HB3	37:DP:100:ARG:NH1	2.14	0.42
37:DP:46:VAL:HB	37:DP:65:ASN:OD1	2.19	0.42
38:DQ:77:LYS:O	38:DQ:80:ASN:HB3	2.19	0.42
24:DV:23:ALA:O	24:DV:24:ASN:HB2	2.19	0.42
44:DX:16:THR:HG23	44:DX:21:LEU:HD12	2.00	0.42
44:DX:37:LEU:HG	44:DX:38:GLN:N	2.35	0.42
1:AA:1096:C:O2'	1:AA:1097:C:H5'	2.20	0.42
1:AA:1111:A:H2	2:AC:176:THR:HG23	1.85	0.42
1:AA:1286:U:O5'	1:AA:1286:U:N1	2.52	0.42
1:AA:1389:C:H2'	1:AA:1390:U:C6	2.55	0.42
1:AA:162:A:H2'	1:AA:163:C:O4'	2.20	0.42
1:AA:545:C:H5''	3:AD:68:GLU:HG2	2.00	0.42
1:AA:806:C:O2'	1:AA:807:A:H5'	2.19	0.42
1:AA:896:C:O2'	1:AA:897:C:H5'	2.19	0.42
20:AB:20:ARG:HG3	20:AB:20:ARG:NH1	2.33	0.42
1:AA:829:G:H4'	20:AB:24:PRO:HG3	2.02	0.42
2:AC:84:GLU:HA	2:AC:87:ARG:NH2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:14:ASP:O	6:AG:18:GLY:HA2	2.19	0.42
15:AP:1:MET:H3	15:AP:24:SER:HB3	1.82	0.42
1:AA:734:G:O2'	17:AR:59:LYS:HD3	2.19	0.42
48:B1:7:LYS:HG3	48:B1:7:LYS:O	2.20	0.42
49:B2:13:ASN:CB	49:B2:18:PHE:HD2	2.29	0.42
23:BB:651:G:C5'	50:B3:17:GLY:HA3	2.50	0.42
22:BA:50:A:H5''	36:BO:68:LYS:CG	2.49	0.42
23:BB:1255:U:C2'	27:BE:67:ARG:HB3	2.50	0.42
23:BB:1322:A:C5	23:BB:1323:C:C5	3.08	0.42
23:BB:1390:U:O2'	23:BB:1391:U:H5'	2.20	0.42
23:BB:159:G:O2'	23:BB:160:A:H5''	2.20	0.42
23:BB:1616:A:H4'	23:BB:1617:C:OP2	2.19	0.42
23:BB:1791:A:C2	23:BB:1829:A:O4'	2.73	0.42
23:BB:1987:A:H2'	23:BB:1988:G:C8	2.54	0.42
23:BB:2019:A:C6	23:BB:2020:A:N7	2.88	0.42
23:BB:2543:G:H2'	23:BB:2544:G:O4'	2.20	0.42
23:BB:2728:U:H5'	32:BK:70:ARG:NH2	2.34	0.42
23:BB:319:G:H2'	23:BB:320:A:C8	2.55	0.42
23:BB:724:U:H2'	23:BB:725:G:O4'	2.19	0.42
23:BB:948:C:H2'	23:BB:949:G:C8	2.53	0.42
25:BC:16:VAL:C	25:BC:17:LYS:HD2	2.40	0.42
25:BC:225:ASN:N	25:BC:226:PRO:HD3	2.35	0.42
25:BC:244:VAL:O	25:BC:245:THR:C	2.58	0.42
25:BC:251:THR:O	25:BC:252:LYS:HB2	2.19	0.42
23:BB:2228:G:OP1	25:BC:257:ARG:HB2	2.20	0.42
26:BD:170:VAL:O	26:BD:171:THR:HB	2.19	0.42
27:BE:43:THR:HG23	27:BE:44:ARG:CZ	2.50	0.42
27:BE:77:ILE:HB	27:BE:78:TRP:CE3	2.54	0.42
30:BH:3:VAL:HG22	30:BH:21:VAL:CG1	2.45	0.42
31:BJ:35:ARG:HH21	31:BJ:40:HIS:H	1.66	0.42
33:BL:27:LEU:C	33:BL:29:LYS:N	2.70	0.42
33:BL:69:ARG:HG3	33:BL:69:ARG:O	2.19	0.42
34:BM:17:ASN:HB3	34:BM:18:ARG:H	1.49	0.42
37:BP:18:SER:HB2	37:BP:87:ARG:CZ	2.50	0.42
23:BB:1156:A:H62	38:BQ:47:ARG:HD3	1.85	0.42
38:BQ:108:LEU:HD22	39:BR:52:PRO:HA	2.00	0.42
42:BU:10:VAL:HG11	42:BU:25:LYS:HZ3	1.84	0.42
42:BU:84:PHE:HD2	42:BU:87:GLU:HB2	1.84	0.42
24:BV:4:ILE:N	24:BV:62:THR:O	2.52	0.42
24:BV:66:ASP:O	24:BV:66:ASP:OD1	2.37	0.42
1:CA:234:C:H2'	1:CA:235:C:H6	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:393:A:O2'	1:CA:394:G:H5'	2.20	0.42
1:CA:489:C:H2'	1:CA:490:C:C6	2.54	0.42
1:CA:925:G:C2	1:CA:927:G:C8	3.08	0.42
1:CA:939:G:H2'	1:CA:940:C:C6	2.54	0.42
2:CC:33:ASP:O	2:CC:37:LYS:HG3	2.19	0.42
2:CC:54:ILE:O	2:CC:54:ILE:HG23	2.20	0.42
3:CD:142:VAL:HG22	3:CD:179:GLY:HA3	2.00	0.42
1:CA:939:G:C5'	6:CG:101:ARG:HH22	2.29	0.42
9:CJ:68:ARG:HD3	9:CJ:70:HIS:NE2	2.35	0.42
11:CL:23:LEU:C	11:CL:25:ALA:N	2.72	0.42
12:CM:1:ALA:O	12:CM:3:ILE:HG13	2.19	0.42
16:CQ:7:LEU:O	16:CQ:59:GLU:HA	2.19	0.42
47:D0:28:SER:HB3	47:D0:36:LYS:O	2.19	0.42
48:D1:9:LYS:HA	48:D1:24:LYS:HB2	2.02	0.42
50:D3:37:THR:HA	50:D3:40:LYS:HD3	2.02	0.42
22:DA:38:C:H2'	22:DA:39:A:O4'	2.18	0.42
22:DA:49:C:C5'	36:DO:101:GLY:HA3	2.50	0.42
23:DB:978:G:O4'	23:DB:1001:A:H2	2.03	0.42
23:DB:1080:A:H4'	52:DI:126:ARG:CD	2.50	0.42
23:DB:1254:A:C6	27:DE:77:ILE:HD12	2.55	0.42
23:DB:1268:A:H2'	23:DB:1269:A:H8	1.85	0.42
23:DB:1410:G:H2'	23:DB:1411:U:H6	1.77	0.42
23:DB:2080:A:H5'	46:DZ:17:SER:HB2	2.00	0.42
23:DB:2183:A:H8	23:DB:2183:A:O5'	2.02	0.42
23:DB:2341:G:H2'	23:DB:2342:C:H6	1.84	0.42
23:DB:2365:G:H4'	43:DW:65:LYS:CD	2.49	0.42
23:DB:2065:C:H1'	23:DB:2449:U:O2	2.20	0.42
23:DB:2840:C:C2	23:DB:2841:C:C5	3.07	0.42
23:DB:340:A:O2'	23:DB:341:C:H5'	2.20	0.42
23:DB:396:G:O2'	23:DB:397:U:H5'	2.20	0.42
23:DB:414:C:H1'	23:DB:1864:U:H1'	2.01	0.42
23:DB:607:U:C4	23:DB:620:G:O4'	2.73	0.42
23:DB:748:G:O5'	40:DS:89:ALA:HB2	2.20	0.42
23:DB:848:C:H2'	23:DB:849:A:H8	1.84	0.42
23:DB:929:U:O2	45:DY:25:GLY:HA2	2.20	0.42
27:DE:150:THR:HG23	27:DE:167:VAL:HG11	2.02	0.42
29:DG:10:VAL:N	29:DG:11:PRO:CD	2.83	0.42
52:DI:100:ILE:HG23	52:DI:104:GLN:OE1	2.20	0.42
33:DL:118:THR:HG21	33:DL:137:ALA:HB3	2.00	0.42
33:DL:14:LYS:CD	33:DL:15:ALA:N	2.83	0.42
23:DB:2394:C:P	33:DL:63:LYS:HD2	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:31:PHE:O	34:DM:102:LEU:HA	2.20	0.42
38:DQ:52:ARG:CB	38:DQ:52:ARG:HH11	2.32	0.42
40:DS:28:LYS:O	40:DS:29:VAL:HB	2.20	0.42
41:DT:31:VAL:HG23	41:DT:83:ALA:O	2.20	0.42
42:DU:64:ILE:HG22	42:DU:64:ILE:O	2.19	0.42
23:DB:2276:G:N7	43:DW:11:ASN:ND2	2.66	0.42
46:DZ:36:VAL:HG12	46:DZ:42:PRO:HB2	2.01	0.42
1:AA:1091:U:H2'	1:AA:1093:A:OP2	2.19	0.42
1:AA:1441:A:H2'	1:AA:1442:G:O4'	2.20	0.42
1:AA:1477:U:H2'	1:AA:1478:U:H6	1.83	0.42
1:AA:21:G:H1'	1:AA:914:A:H61	1.84	0.42
1:AA:257:G:H2'	1:AA:258:G:C8	2.55	0.42
1:AA:489:C:H2'	1:AA:490:C:C6	2.55	0.42
1:AA:309:A:O2'	1:AA:607:A:N1	2.52	0.42
1:AA:771:G:H2'	1:AA:772:U:C6	2.54	0.42
1:AA:1073:U:O3'	20:AB:104:LYS:NZ	2.53	0.42
20:AB:53:LEU:HB3	20:AB:219:THR:HG21	2.02	0.42
20:AB:80:LYS:HB3	20:AB:90:PHE:CE1	2.54	0.42
2:AC:181:ILE:HG22	2:AC:182:ASP:N	2.35	0.42
2:AC:38:VAL:HG12	2:AC:93:ILE:HG21	2.01	0.42
3:AD:119:HIS:C	3:AD:121:ALA:H	2.22	0.42
4:AE:110:MET:SD	4:AE:110:MET:N	2.92	0.42
4:AE:50:GLY:H	4:AE:62:ALA:HB2	1.85	0.42
5:AF:77:THR:O	5:AF:81:ASN:HB2	2.20	0.42
6:AG:148:LYS:HA	6:AG:151:ALA:HB3	2.02	0.42
6:AG:30:MET:SD	6:AG:35:LYS:HB2	2.59	0.42
8:AI:24:ASN:HA	8:AI:26:LYS:HZ2	1.84	0.42
12:AM:14:ALA:CB	12:AM:33:LEU:HD11	2.50	0.42
14:AO:10:ILE:HD13	14:AO:30:LEU:HA	2.02	0.42
15:AP:71:VAL:HA	15:AP:74:LEU:CD1	2.49	0.42
16:AQ:26:ARG:O	16:AQ:38:LYS:HA	2.19	0.42
16:AQ:62:GLU:HG3	16:AQ:72:TRP:CH2	2.55	0.42
18:AS:20:LYS:NZ	18:AS:20:LYS:HB3	2.31	0.42
18:AS:47:THR:HA	18:AS:59:VAL:O	2.20	0.42
12:AM:84:CYS:HA	18:AS:72:GLU:O	2.19	0.42
21:AU:45:LYS:O	21:AU:48:LYS:HG3	2.19	0.42
48:B1:34:GLU:N	48:B1:52:LYS:O	2.53	0.42
23:BB:100:U:O2'	23:BB:101:A:O5'	2.38	0.42
23:BB:1268:A:H2'	23:BB:1269:A:H8	1.85	0.42
23:BB:1528:A:C2'	23:BB:1529:G:H5'	2.50	0.42
23:BB:1797:G:C5'	25:BC:251:THR:HG23	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1820:U:H3	25:BC:197:ALA:CB	2.33	0.42
23:BB:18:U:O3'	38:BQ:22:GLY:HA3	2.19	0.42
23:BB:2054:A:OP1	23:BB:2055:C:H4'	2.20	0.42
23:BB:2299:U:H2'	23:BB:2300:C:H6	1.85	0.42
23:BB:2428:G:H5''	23:BB:2429:G:O5'	2.19	0.42
23:BB:2733:A:O5'	23:BB:2733:A:H8	2.03	0.42
23:BB:2733:A:N1	26:BD:209:ALA:O	2.53	0.42
23:BB:288:U:H2'	23:BB:289:G:H8	1.85	0.42
23:BB:46:G:H2'	23:BB:47:C:C6	2.55	0.42
23:BB:45:G:H5'	23:BB:46:G:OP1	2.19	0.42
23:BB:729:G:H8	23:BB:764:A:OP1	2.03	0.42
23:BB:834:G:H2'	23:BB:835:C:C6	2.54	0.42
22:BA:100:G:H5''	23:BB:915:C:O2'	2.19	0.42
25:BC:159:THR:C	25:BC:160:TYR:HD2	2.23	0.42
25:BC:191:LEU:O	25:BC:191:LEU:HD13	2.20	0.42
25:BC:73:ILE:N	25:BC:73:ILE:HD12	2.35	0.42
26:BD:29:VAL:HG23	26:BD:51:THR:CG2	2.50	0.42
27:BE:145:ASP:C	27:BE:147:LEU:N	2.71	0.42
30:BH:6:LEU:H	30:BH:16:GLY:N	2.17	0.42
31:BJ:120:ARG:HD3	31:BJ:120:ARG:O	2.20	0.42
32:BK:119:ALA:N	32:BK:120:PRO:CD	2.82	0.42
32:BK:24:VAL:HG22	32:BK:39:ILE:HD12	2.01	0.42
33:BL:94:THR:HG23	33:BL:95:LEU:N	2.34	0.42
34:BM:43:ALA:C	34:BM:45:GLN:H	2.23	0.42
35:BN:36:THR:O	35:BN:37:THR:O	2.37	0.42
37:BP:92:ARG:H	37:BP:113:LEU:HD23	1.84	0.42
38:BQ:58:GLN:O	38:BQ:62:ALA:HB2	2.19	0.42
41:BT:6:ARG:HB2	41:BT:6:ARG:HE	1.57	0.42
42:BU:25:LYS:HA	42:BU:35:VAL:N	2.29	0.42
42:BU:62:ALA:O	42:BU:63:ALA:C	2.58	0.42
24:BV:44:HIS:CE1	24:BV:85:LYS:HB2	2.55	0.42
43:BW:54:ARG:CZ	43:BW:54:ARG:HB2	2.50	0.42
44:BX:29:ARG:HG2	44:BX:29:ARG:O	2.20	0.42
1:CA:257:G:H2'	1:CA:258:G:C8	2.55	0.42
1:CA:297:G:N2	1:CA:300:A:OP2	2.51	0.42
1:CA:462:G:H3'	1:CA:463:U:C5'	2.48	0.42
1:CA:687:A:C2	1:CA:704:A:C6	3.08	0.42
1:CA:761:G:H2'	1:CA:762:U:H6	1.84	0.42
1:CA:771:G:H2'	1:CA:772:U:C6	2.54	0.42
20:CB:163:ILE:CG2	20:CB:164:ASP:H	2.25	0.42
20:CB:202:ASN:HD22	20:CB:203:ASP:N	2.12	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CC:68:HIS:HA	2:CC:103:ALA:HB3	2.02	0.42
3:CD:55:ARG:HD2	3:CD:55:ARG:HA	1.78	0.42
3:CD:67:LEU:O	3:CD:71:PHE:HB2	2.19	0.42
4:CE:144:GLU:HA	4:CE:146:MET:SD	2.59	0.42
6:CG:11:ILE:HG22	6:CG:12:LEU:H	1.84	0.42
1:CA:716:A:N3	10:CK:119:GLY:HA2	2.34	0.42
10:CK:74:LYS:C	10:CK:76:TYR:H	2.23	0.42
2:CC:5:HIS:ND1	13:CN:88:MET:HB3	2.35	0.42
14:CO:10:ILE:HD13	14:CO:30:LEU:HG	2.02	0.42
15:CP:4:ILE:HG22	15:CP:19:VAL:HG23	2.02	0.42
15:CP:6:LEU:HB3	15:CP:17:TYR:HB3	2.01	0.42
16:CQ:82:VAL:C	16:CQ:83:LEU:HD22	2.40	0.42
19:CT:67:HIS:O	19:CT:70:LYS:HG2	2.19	0.42
47:D0:50:GLY:O	47:D0:51:ARG:C	2.57	0.42
23:DB:1161:C:H2'	23:DB:1162:G:C8	2.53	0.42
23:DB:1289:C:C2	23:DB:1290:C:C5	3.07	0.42
23:DB:1463:C:H2'	23:DB:1464:G:H8	1.82	0.42
23:DB:2077:A:N6	23:DB:2435:A:N6	2.68	0.42
23:DB:2617:U:C4	23:DB:2618:G:N7	2.87	0.42
23:DB:265:A:H2'	23:DB:266:G:C1'	2.50	0.42
23:DB:357:C:H2'	23:DB:358:U:C6	2.54	0.42
23:DB:587:C:H5''	33:DL:29:LYS:HZ2	1.84	0.42
23:DB:705:A:N6	23:DB:726:G:H1'	2.35	0.42
23:DB:795:C:H2'	23:DB:796:C:H6	1.84	0.42
23:DB:822:G:H2'	23:DB:823:C:H6	1.84	0.42
23:DB:86:G:P	42:DU:28:LEU:HD22	2.60	0.42
23:DB:876:C:H5'	23:DB:877:A:OP2	2.20	0.42
23:DB:963:U:H5''	55:DB:3696:HOH:O	2.19	0.42
25:DC:179:GLU:OE2	25:DC:266:ILE:HA	2.19	0.42
26:DD:118:PHE:CA	26:DD:164:GLN:HG2	2.50	0.42
26:DD:122:VAL:HB	26:DD:141:ARG:NH1	2.35	0.42
26:DD:156:PHE:CG	31:DJ:81:ILE:HG21	2.54	0.42
26:DD:199:SER:O	26:DD:201:LEU:HD13	2.20	0.42
26:DD:77:ARG:NH2	26:DD:79:LEU:HB2	2.14	0.42
29:DG:3:VAL:CG2	29:DG:4:ALA:N	2.83	0.42
29:DG:29:ASN:HB2	29:DG:78:VAL:HA	2.01	0.42
30:DH:8:LYS:HD2	30:DH:14:SER:HA	2.02	0.42
52:DI:63:ASP:O	52:DI:63:ASP:OD1	2.38	0.42
34:DM:33:LEU:H	34:DM:101:VAL:HB	1.85	0.42
34:DM:127:LYS:NZ	34:DM:129:THR:N	2.68	0.42
34:DM:38:ARG:HD3	34:DM:39:GLY:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DM:96:ILE:HG22	34:DM:97:GLN:CD	2.39	0.42
35:DN:34:ILE:HG22	35:DN:35:LYS:N	2.28	0.42
35:DN:7:GLY:O	35:DN:8:ARG:HB2	2.20	0.42
37:DP:73:PHE:CD2	37:DP:75:THR:HG23	2.55	0.42
38:DQ:65:ASN:CB	38:DQ:75:TYR:HB2	2.49	0.42
39:DR:22:LEU:CD1	39:DR:23:GLU:H	2.28	0.42
39:DR:46:GLU:HB3	39:DR:47:VAL:H	1.63	0.42
40:DS:68:ASP:C	40:DS:69:LEU:HD22	2.40	0.42
41:DT:66:LYS:HB3	41:DT:66:LYS:NZ	2.35	0.42
42:DU:49:PRO:CD	42:DU:50:ALA:H	2.30	0.42
44:DX:31:GLN:HE21	44:DX:31:GLN:CA	2.24	0.42
1:AA:109:A:N1	1:AA:326:G:C6	2.88	0.42
1:AA:1110:A:C5	1:AA:1111:A:N7	2.88	0.42
1:AA:113:G:H21	1:AA:353:A:H8	1.66	0.42
1:AA:538:G:H2'	1:AA:539:A:H8	1.85	0.42
1:AA:6:G:H4'	1:AA:298:A:H4'	2.01	0.42
1:AA:839:C:O2'	1:AA:840:C:H5'	2.19	0.42
1:AA:811:C:O2'	1:AA:901:A:N1	2.47	0.42
1:AA:959:A:H2'	1:AA:960:U:O4'	2.20	0.42
1:AA:960:U:O2'	1:AA:1223:C:H4'	2.20	0.42
20:AB:107:ARG:HA	20:AB:110:ILE:HD12	2.01	0.42
20:AB:65:LYS:HB2	20:AB:157:PRO:HA	2.02	0.42
2:AC:57:GLU:O	2:AC:63:ILE:HA	2.20	0.42
3:AD:137:SER:HB2	3:AD:138:PRO:CD	2.45	0.42
3:AD:54:LEU:HA	3:AD:57:LYS:HB3	2.02	0.42
5:AF:70:VAL:HA	5:AF:73:GLU:HG3	2.02	0.42
4:AE:82:HIS:CD2	7:AH:95:MET:HG3	2.55	0.42
9:AJ:10:LEU:O	9:AJ:71:LEU:HA	2.19	0.42
11:AL:41:PRO:HG3	11:AL:46:SER:O	2.19	0.42
12:AM:106:ARG:HD3	12:AM:106:ARG:HA	1.90	0.42
13:AN:1:ALA:O	13:AN:5:MET:HB2	2.19	0.42
2:AC:5:HIS:CG	13:AN:88:MET:HB3	2.55	0.42
16:AQ:9:GLY:O	16:AQ:58:VAL:HG22	2.20	0.42
47:B0:49:ARG:O	47:B0:49:ARG:HG2	2.20	0.42
48:B1:46:VAL:C	48:B1:47:ILE:HG12	2.41	0.42
49:B2:34:ARG:HG3	49:B2:34:ARG:NH1	2.35	0.42
50:B3:12:ARG:HG2	50:B3:23:HIS:CB	2.50	0.42
23:BB:1030:C:O2'	23:BB:1031:G:H5'	2.20	0.42
23:BB:1557:C:H3'	23:BB:1558:C:C5'	2.49	0.42
23:BB:1700:A:C2'	23:BB:1701:A:H5'	2.44	0.42
1:AA:1473:G:O2'	23:BB:1702:G:H4'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:1799:G:O2'	25:BC:181:ARG:HD2	2.19	0.42
23:BB:1829:A:H3'	23:BB:1830:C:C6	2.55	0.42
23:BB:1897:G:O2'	23:BB:1898:U:H5'	2.20	0.42
23:BB:2078:C:H2'	23:BB:2079:U:C6	2.55	0.42
23:BB:2328:A:O2'	23:BB:2329:U:H5'	2.19	0.42
23:BB:2351:G:HO2'	23:BB:2352:A:H8	1.68	0.42
23:BB:1783:A:H5'	23:BB:2608:G:H4'	2.02	0.42
23:BB:2684:U:OP1	37:BP:59:THR:HG22	2.19	0.42
23:BB:2756:U:H5'	51:B4:18:LYS:HZ1	1.85	0.42
23:BB:2849:U:O4'	23:BB:2868:A:C6	2.73	0.42
23:BB:39:G:H2'	23:BB:40:U:H6	1.84	0.42
23:BB:953:G:H2'	23:BB:954:G:H8	1.85	0.42
25:BC:161:VAL:HB	25:BC:193:GLU:OE2	2.20	0.42
25:BC:242:HIS:N	25:BC:243:PRO:HD3	2.34	0.42
25:BC:180:MET:HB2	25:BC:265:PHE:CB	2.49	0.42
25:BC:48:ILE:HG22	25:BC:49:THR:N	2.35	0.42
26:BD:139:SER:O	26:BD:140:HIS:C	2.58	0.42
26:BD:38:LYS:O	26:BD:42:ASN:ND2	2.49	0.42
26:BD:2:ILE:HG23	26:BD:3:GLY:N	2.33	0.42
27:BE:152:GLU:H	27:BE:171:ASP:CB	2.33	0.42
28:BF:56:LEU:HD23	28:BF:59:ILE:CG2	2.50	0.42
29:BG:132:LEU:HD21	29:BG:147:LEU:CD1	2.49	0.42
29:BG:28:LYS:C	29:BG:29:ASN:HD22	2.23	0.42
29:BG:66:THR:O	29:BG:70:LEU:HB2	2.20	0.42
30:BH:142:VAL:HG12	30:BH:143:ILE:H	1.84	0.42
30:BH:82:SER:HB3	30:BH:90:LEU:HD23	2.02	0.42
52:BI:12:VAL:HG23	52:BI:41:PHE:CZ	2.55	0.42
52:BI:64:ARG:C	52:BI:64:ARG:HD2	2.39	0.42
52:BI:5:GLN:CG	52:BI:6:ALA:N	2.81	0.42
33:BL:38:GLN:N	33:BL:41:ARG:NH1	2.68	0.42
34:BM:3:GLN:HG2	34:BM:46:ILE:CB	2.44	0.42
34:BM:75:GLU:O	34:BM:86:LYS:HB2	2.20	0.42
35:BN:72:ASP:C	35:BN:74:GLU:H	2.24	0.42
36:BO:55:GLU:HG3	36:BO:56:LYS:H	1.84	0.42
36:BO:6:ALA:C	36:BO:8:ILE:H	2.23	0.42
37:BP:80:VAL:HG22	37:BP:80:VAL:O	2.20	0.42
37:BP:91:VAL:HG12	37:BP:92:ARG:N	2.35	0.42
41:BT:84:TYR:O	41:BT:85:VAL:O	2.38	0.42
43:BW:44:PHE:N	43:BW:76:ARG:HB2	2.35	0.42
1:CA:1053:G:H4'	1:CA:1054:C:H5'	2.00	0.42
1:CA:123:U:H2'	1:CA:124:C:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1381:U:O2'	1:CA:1382:C:H5'	2.20	0.42
1:CA:1499:A:H1'	1:CA:1520:C:H5'	2.01	0.42
1:CA:415:A:N3	1:CA:415:A:O4'	2.53	0.42
1:CA:635:A:H2'	1:CA:636:U:H6	1.85	0.42
1:CA:688:G:O2'	1:CA:689:C:H5'	2.20	0.42
1:CA:989:U:O2'	1:CA:990:C:H5'	2.20	0.42
20:CB:209:VAL:HG23	20:CB:210:THR:H	1.85	0.42
20:CB:27:LYS:CB	20:CB:28:PRO:HD3	2.46	0.42
2:CC:163:ARG:O	2:CC:164:THR:HG23	2.20	0.42
2:CC:28:PHE:CZ	13:CN:93:PRO:HG2	2.55	0.42
2:CC:37:LYS:HB3	2:CC:41:TYR:CZ	2.55	0.42
4:CE:17:VAL:HG22	4:CE:17:VAL:O	2.20	0.42
5:CF:29:ILE:HG23	5:CF:66:ALA:HB2	2.02	0.42
6:CG:55:LYS:CG	6:CG:56:SER:H	2.30	0.42
6:CG:56:SER:HB3	6:CG:59:GLU:CB	2.50	0.42
9:CJ:9:ARG:HG3	9:CJ:99:GLN:HB2	2.01	0.42
13:CN:26:LEU:HA	13:CN:29:ILE:CD1	2.44	0.42
13:CN:30:ILE:H	13:CN:44:VAL:HG21	1.85	0.42
14:CO:69:LEU:HD21	14:CO:76:ARG:HD2	2.02	0.42
15:CP:12:LYS:HD2	15:CP:13:LYS:HG3	2.02	0.42
15:CP:67:ILE:HG13	15:CP:71:VAL:CG1	2.40	0.42
1:CA:280:C:N3	16:CQ:39:ARG:HA	2.35	0.42
18:CS:46:LEU:HB3	18:CS:47:THR:H	1.73	0.42
18:CS:30:LEU:HB2	18:CS:48:ILE:HG12	2.01	0.42
19:CT:47:GLN:HG2	19:CT:82:ILE:HD12	2.01	0.42
47:D0:42:ILE:O	47:D0:43:THR:C	2.57	0.42
50:D3:46:LYS:HD2	50:D3:47:ALA:N	2.35	0.42
23:DB:1099:G:C8	52:DI:3:LYS:O	2.73	0.42
23:DB:1399:C:O2'	23:DB:1400:U:H5'	2.20	0.42
23:DB:161:A:H3'	23:DB:162:U:C5'	2.38	0.42
23:DB:162:U:O2'	23:DB:163:C:C5'	2.60	0.42
23:DB:1665:A:H1'	32:DK:67:LYS:HE2	2.02	0.42
23:DB:1674:G:N2	23:DB:1677:A:N1	2.67	0.42
23:DB:1713:A:N6	23:DB:1745:A:H61	2.18	0.42
23:DB:1791:A:C2	23:DB:1829:A:O4'	2.73	0.42
23:DB:211:C:O2'	23:DB:212:G:H5'	2.20	0.42
23:DB:2266:A:O4'	23:DB:2272:U:O4	2.37	0.42
23:DB:2347:C:H4'	23:DB:2347:C:OP1	2.19	0.42
23:DB:2527:C:H4'	51:D4:34:LYS:O	2.20	0.42
23:DB:409:G:H2'	23:DB:410:G:H8	1.85	0.42
23:DB:459:U:O2'	23:DB:460:A:H5'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:647:G:H2'	23:DB:648:G:H8	1.85	0.42
25:DC:91:ALA:O	25:DC:102:TYR:HD2	2.03	0.42
25:DC:103:ILE:CG2	25:DC:104:LEU:H	2.33	0.42
25:DC:161:VAL:O	25:DC:161:VAL:HG12	2.19	0.42
25:DC:257:ARG:HA	25:DC:261:ARG:CZ	2.49	0.42
26:DD:32:ASN:O	26:DD:34:VAL:HG13	2.20	0.42
26:DD:34:VAL:HG11	26:DD:50:VAL:CG2	2.50	0.42
27:DE:169:VAL:O	27:DE:170:ARG:HB2	2.19	0.42
27:DE:189:THR:C	27:DE:191:ASP:N	2.72	0.42
28:DF:163:GLU:HA	28:DF:166:ARG:CG	2.50	0.42
29:DG:155:PRO:O	29:DG:171:LYS:HG3	2.19	0.42
29:DG:29:ASN:OD1	29:DG:81:GLY:HA2	2.19	0.42
30:DH:101:ASP:O	30:DH:104:THR:HB	2.19	0.42
30:DH:131:SER:HB2	30:DH:141:LYS:HG3	2.02	0.42
30:DH:144:VAL:HG12	30:DH:145:ASN:N	2.34	0.42
30:DH:35:LYS:O	30:DH:35:LYS:HG3	2.20	0.42
52:DI:45:THR:C	52:DI:48:ILE:HG22	2.40	0.42
31:DJ:44:TYR:CE2	31:DJ:50:THR:HB	2.54	0.42
31:DJ:67:ASN:C	31:DJ:69:ARG:N	2.73	0.42
31:DJ:96:ARG:HG2	31:DJ:99:ARG:HB2	2.01	0.42
34:DM:16:ARG:HH11	34:DM:16:ARG:CG	2.33	0.42
35:DN:12:ARG:HB3	35:DN:13:ASN:H	1.77	0.42
37:DP:26:GLU:O	37:DP:27:VAL:C	2.59	0.42
37:DP:18:SER:HA	37:DP:87:ARG:HH22	1.85	0.42
38:DQ:92:LYS:C	38:DQ:94:LEU:N	2.73	0.42
38:DQ:92:LYS:O	38:DQ:94:LEU:N	2.53	0.42
41:DT:53:VAL:HB	41:DT:87:LEU:HD21	2.02	0.42
42:DU:40:LEU:O	42:DU:58:VAL:HA	2.19	0.42
43:DW:75:ASN:O	43:DW:77:LYS:N	2.52	0.42
44:DX:55:THR:O	44:DX:57:LEU:N	2.40	0.42
46:DZ:59:ARG:O	46:DZ:60:PHE:HB3	2.20	0.42
1:AA:1052:U:H5'	1:AA:1053:G:OP2	2.20	0.41
1:AA:207:C:O2'	1:AA:208:U:H5'	2.20	0.41
1:AA:208:U:C2'	1:AA:209:U:H5''	2.50	0.41
1:AA:412:A:H1'	1:AA:413:G:C8	2.46	0.41
1:AA:821:G:H2'	1:AA:822:U:C6	2.55	0.41
1:AA:952:U:O2'	1:AA:953:G:H5'	2.20	0.41
2:AC:172:VAL:HG11	2:AC:200:TRP:HB3	2.02	0.41
2:AC:24:ASN:O	2:AC:26:LYS:N	2.54	0.41
4:AE:137:ARG:HG2	4:AE:137:ARG:H	1.75	0.41
4:AE:82:HIS:HE1	4:AE:147:ASN:H	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:15:ALA:O	8:AI:66:VAL:HA	2.20	0.41
8:AI:93:LEU:C	8:AI:93:LEU:HD12	2.40	0.41
9:AJ:93:ALA:HB1	9:AJ:96:VAL:HG23	2.02	0.41
10:AK:105:ARG:HB3	10:AK:106:ILE:H	1.65	0.41
10:AK:121:ARG:HG3	10:AK:121:ARG:NH1	2.34	0.41
10:AK:22:ILE:HG23	10:AK:31:VAL:HG22	2.02	0.41
13:AN:19:TYR:HE2	13:AN:23:ARG:NH1	2.17	0.41
13:AN:20:PHE:HA	13:AN:24:ALA:H	1.85	0.41
14:AO:52:ARG:O	14:AO:55:LEU:HB3	2.20	0.41
15:AP:8:ARG:NE	15:AP:15:PRO:HB3	2.34	0.41
15:AP:51:ARG:O	15:AP:52:LEU:HD13	2.19	0.41
15:AP:68:SER:HB3	15:AP:71:VAL:CG1	2.47	0.41
18:AS:44:ILE:HA	18:AS:61:VAL:HB	2.01	0.41
47:B0:15:ARG:HG2	47:B0:15:ARG:NH1	2.35	0.41
47:B0:4:GLN:HB3	47:B0:4:GLN:HE21	1.65	0.41
48:B1:49:LYS:H	48:B1:49:LYS:HD2	1.85	0.41
50:B3:24:LYS:HE2	50:B3:29:ARG:HH22	1.85	0.41
23:BB:1082:U:H2'	23:BB:1083:U:O4'	2.20	0.41
23:BB:1236:G:O2'	23:BB:1237:A:O4'	2.36	0.41
23:BB:1268:A:H2'	23:BB:1269:A:C8	2.55	0.41
23:BB:1289:C:C2	23:BB:1290:C:C5	3.07	0.41
23:BB:1294:U:C4	23:BB:1295:C:C5	3.07	0.41
23:BB:136:G:P	23:BB:136:G:H8	2.43	0.41
23:BB:1372:U:H2'	23:BB:1373:A:H8	1.86	0.41
23:BB:1529:G:H2'	23:BB:1530:G:H8	1.85	0.41
23:BB:161:A:O5'	23:BB:161:A:H8	2.02	0.41
23:BB:1678:A:C2'	23:BB:1679:A:H5'	2.50	0.41
23:BB:2026:U:C2	23:BB:2027:G:C8	3.07	0.41
23:BB:2046:G:H5'	47:B0:15:ARG:HG3	2.01	0.41
23:BB:2047:C:OP1	47:B0:15:ARG:HD2	2.20	0.41
23:BB:2308:G:H5'	23:BB:2309:A:C5'	2.49	0.41
23:BB:2322:A:C6	23:BB:2333:A:N6	2.88	0.41
23:BB:2324:U:O2'	23:BB:2385:C:H5	2.02	0.41
23:BB:2708:G:O2'	23:BB:2709:G:H5'	2.20	0.41
23:BB:354:A:H2'	23:BB:355:U:O4'	2.20	0.41
23:BB:433:C:H2'	23:BB:434:U:H6	1.84	0.41
23:BB:482:A:H5''	42:BU:43:LYS:CE	2.33	0.41
23:BB:508:A:H2'	23:BB:509:C:C5'	2.45	0.41
23:BB:522:A:H2'	23:BB:523:C:H6	1.85	0.41
23:BB:58:G:H2'	23:BB:59:U:C6	2.55	0.41
23:BB:647:G:H2'	23:BB:648:G:H8	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:908:C:O2'	23:BB:909:A:H5'	2.20	0.41
25:BC:51:ARG:O	25:BC:51:ARG:NE	2.52	0.41
26:BD:100:LEU:HG	26:BD:100:LEU:O	2.20	0.41
27:BE:18:THR:O	27:BE:19:PHE:C	2.58	0.41
27:BE:84:THR:CG2	27:BE:85:PHE:H	2.26	0.41
30:BH:98:ASP:O	30:BH:102:ALA:HB2	2.20	0.41
30:BH:89:LYS:HE2	30:BH:90:LEU:N	2.35	0.41
32:BK:2:ILE:CD1	32:BK:8:LEU:HD21	2.50	0.41
32:BK:87:LEU:O	32:BK:89:ASN:N	2.53	0.41
23:BB:636:G:N1	33:BL:76:GLU:OE2	2.53	0.41
34:BM:10:ARG:HB3	34:BM:11:LYS:HE3	2.02	0.41
35:BN:65:LEU:HD12	35:BN:68:ALA:HB3	2.02	0.41
22:BA:50:A:OP1	36:BO:68:LYS:HG3	2.19	0.41
37:BP:52:ARG:HB3	37:BP:60:VAL:HG21	2.00	0.41
38:BQ:65:ASN:HB2	38:BQ:75:TYR:CB	2.45	0.41
38:BQ:94:LEU:H	38:BQ:94:LEU:CD1	2.27	0.41
42:BU:92:VAL:HG22	42:BU:92:VAL:O	2.20	0.41
22:BA:75:G:H1'	24:BV:29:ILE:HG21	2.01	0.41
24:BV:76:ASP:HB2	34:BM:136:MET:CE	2.49	0.41
23:BB:2270:A:O5'	43:BW:18:LYS:HB3	2.19	0.41
43:BW:40:ARG:NH1	43:BW:68:PHE:HD1	2.17	0.41
45:BY:28:LEU:N	45:BY:28:LEU:HD12	2.35	0.41
1:CA:1279:G:N1	9:CJ:45:ARG:NH1	2.68	0.41
1:CA:1292:G:O2'	1:CA:1293:C:H5'	2.20	0.41
1:CA:1307:U:H2'	1:CA:1308:U:O4'	2.19	0.41
1:CA:154:U:O2'	1:CA:155:A:H5'	2.20	0.41
1:CA:73:C:H2'	1:CA:74:A:H8	1.85	0.41
1:CA:928:G:H2'	1:CA:929:G:H8	1.85	0.41
1:CA:951:G:O2'	1:CA:952:U:H5'	2.19	0.41
1:CA:975:A:O2'	1:CA:976:G:P	2.77	0.41
20:CB:181:PRO:HA	20:CB:196:ASP:OD2	2.19	0.41
2:CC:53:ARG:O	2:CC:67:ILE:HD12	2.20	0.41
2:CC:63:ILE:O	2:CC:63:ILE:HG23	2.19	0.41
3:CD:169:TRP:HB2	3:CD:183:ARG:O	2.20	0.41
3:CD:24:VAL:HG23	3:CD:25:ARG:N	2.35	0.41
3:CD:97:LEU:HB2	3:CD:134:TYR:HB3	2.02	0.41
5:CF:11:HIS:NE2	5:CF:13:ASP:HB3	2.34	0.41
5:CF:85:ILE:CG2	5:CF:86:ARG:H	2.32	0.41
8:CI:105:ARG:O	8:CI:107:ALA:N	2.53	0.41
10:CK:30:ILE:HD12	10:CK:30:ILE:C	2.40	0.41
10:CK:94:SER:HA	10:CK:97:ARG:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:CN:51:PRO:HG2	13:CN:52:ARG:N	2.30	0.41
9:CJ:65:TYR:CE1	13:CN:84:ARG:HA	2.54	0.41
15:CP:1:MET:HA	15:CP:1:MET:HE3	2.01	0.41
16:CQ:45:VAL:HA	16:CQ:72:TRP:O	2.20	0.41
18:CS:10:ILE:HG23	18:CS:40:PHE:HE2	1.85	0.41
1:CA:262:A:H5''	19:CT:70:LYS:HD2	2.02	0.41
47:D0:12:ARG:HG3	47:D0:13:GLY:N	2.35	0.41
48:D1:8:ILE:HB	48:D1:27:ARG:NH1	2.35	0.41
23:DB:1112:G:H5'	29:DG:1:SER:HA	2.01	0.41
23:DB:1221:C:O2'	23:DB:1222:U:H5'	2.20	0.41
23:DB:1373:A:OP1	23:DB:2213:U:C4	2.73	0.41
23:DB:1421:G:C2	23:DB:1422:G:C8	3.08	0.41
23:DB:1529:G:H2'	23:DB:1530:G:H8	1.85	0.41
23:DB:1541:C:H2'	23:DB:1542:U:H6	1.85	0.41
23:DB:1714:U:HO2'	23:DB:1715:G:P	2.42	0.41
23:DB:1788:C:O2'	23:DB:1789:A:H5'	2.20	0.41
23:DB:1812:U:H2'	23:DB:1813:G:C8	2.53	0.41
23:DB:2091:C:H1'	46:DZ:32:LEU:HG	2.02	0.41
23:DB:2708:G:O2'	23:DB:2709:G:H5'	2.20	0.41
23:DB:327:G:H2'	23:DB:328:U:O4'	2.20	0.41
23:DB:35:G:H2'	23:DB:36:G:O4'	2.20	0.41
23:DB:608:A:H2'	23:DB:609:A:H8	1.81	0.41
23:DB:775:G:C5	23:DB:794:A:C8	3.08	0.41
25:DC:22:GLU:HA	25:DC:202:ARG:CZ	2.50	0.41
28:DF:132:ARG:HG3	28:DF:132:ARG:HH11	1.85	0.41
30:DH:3:VAL:CG2	30:DH:37:VAL:HG21	2.49	0.41
31:DJ:41:LYS:O	38:DQ:63:ARG:NH2	2.53	0.41
32:DK:63:VAL:CG2	32:DK:115:ILE:HD11	2.50	0.41
36:DO:10:ARG:HG3	36:DO:96:GLY:N	2.35	0.41
39:DR:32:THR:HA	39:DR:35:PHE:CE2	2.55	0.41
40:DS:8:ARG:HA	40:DS:102:HIS:ND1	2.35	0.41
23:DB:397:U:OP1	46:DZ:48:GLN:HG3	2.20	0.41
1:AA:1414:U:H2'	1:AA:1415:G:C8	2.55	0.41
1:AA:182:A:O2'	1:AA:183:C:O5'	2.34	0.41
1:AA:372:C:C1'	1:AA:373:A:OP2	2.66	0.41
1:AA:598:U:C2	1:AA:599:C:C5	3.09	0.41
1:AA:981:U:P	13:AN:12:ARG:HH12	2.42	0.41
20:AB:185:ILE:HD12	20:AB:185:ILE:O	2.20	0.41
20:AB:14:HIS:NE2	20:AB:49:PHE:HZ	2.17	0.41
2:AC:131:ARG:HB3	2:AC:131:ARG:CZ	2.50	0.41
2:AC:186:SER:HB3	2:AC:197:VAL:CG1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AC:82:ASP:O	2:AC:86:LEU:HG	2.19	0.41
3:AD:43:ARG:HB3	3:AD:44:LYS:H	1.48	0.41
4:AE:61:LYS:O	4:AE:65:LYS:HG2	2.20	0.41
4:AE:63:MET:O	4:AE:66:ALA:HB3	2.20	0.41
6:AG:96:ASN:O	6:AG:100:MET:HG3	2.20	0.41
14:AO:80:LEU:O	14:AO:84:LEU:HD13	2.20	0.41
17:AR:32:ILE:HG13	17:AR:33:THR:N	2.35	0.41
17:AR:70:THR:OG1	17:AR:71:ASP:N	2.52	0.41
19:AT:60:GLN:C	19:AT:62:ALA:H	2.22	0.41
51:B4:24:ARG:HG2	51:B4:26:ILE:HG12	2.02	0.41
22:BA:92:C:H2'	22:BA:93:C:H6	1.84	0.41
23:BB:1010:A:O2'	23:BB:1152:C:H1'	2.20	0.41
23:BB:1119:U:H2'	23:BB:1120:G:H8	1.85	0.41
23:BB:1215:G:O2'	23:BB:1216:G:H5'	2.20	0.41
23:BB:1998:A:H2'	23:BB:1999:C:C6	2.55	0.41
23:BB:2149:U:OP1	23:BB:2149:U:H4'	2.20	0.41
23:BB:2185:U:H6	23:BB:2185:U:O5'	2.02	0.41
23:BB:2462:C:H2'	23:BB:2463:C:H6	1.84	0.41
23:BB:2513:A:H2'	23:BB:2514:U:C6	2.55	0.41
23:BB:2526:G:N2	51:B4:1:MET:HB3	2.35	0.41
23:BB:2673:G:H2'	23:BB:2674:G:H8	1.85	0.41
23:BB:2768:U:H2'	23:BB:2769:U:O4'	2.20	0.41
23:BB:2849:U:N3	23:BB:2867:G:H1'	2.35	0.41
23:BB:30:G:H2'	23:BB:31:C:H6	1.85	0.41
25:BC:163:ILE:CG2	25:BC:164:VAL:H	2.17	0.41
25:BC:251:THR:HG22	25:BC:252:LYS:N	2.34	0.41
25:BC:27:LYS:N	25:BC:28:PRO:CD	2.83	0.41
25:BC:78:GLU:HB3	25:BC:92:LEU:HD23	2.02	0.41
23:BB:2572:A:C2	26:BD:150:GLN:NE2	2.88	0.41
26:BD:33:ARG:O	26:BD:34:VAL:HB	2.19	0.41
52:BI:35:MET:C	52:BI:35:MET:SD	2.98	0.41
33:BL:47:ARG:HG3	33:BL:47:ARG:O	2.20	0.41
33:BL:73:ILE:O	33:BL:74:THR:C	2.59	0.41
23:BB:2723:C:H5''	35:BN:3:HIS:HB2	2.03	0.41
36:BO:39:VAL:HB	36:BO:83:LEU:CD1	2.49	0.41
36:BO:45:SER:C	36:BO:47:VAL:N	2.73	0.41
36:BO:49:VAL:CG2	36:BO:50:ALA:H	2.19	0.41
40:BS:8:ARG:NH1	40:BS:102:HIS:NE2	2.67	0.41
41:BT:16:VAL:HG13	41:BT:16:VAL:O	2.20	0.41
42:BU:35:VAL:CG1	42:BU:36:GLU:H	2.15	0.41
42:BU:71:ILE:HG23	42:BU:72:PHE:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:BV:6:ALA:HB3	24:BV:65:VAL:CG1	2.50	0.41
43:BW:35:ILE:HG12	43:BW:70:VAL:CG2	2.47	0.41
44:BX:19:LEU:C	44:BX:21:LEU:N	2.74	0.41
1:CA:1148:U:H2'	1:CA:1149:C:O4'	2.19	0.41
1:CA:1320:C:H41	18:CS:36:ARG:CG	2.32	0.41
1:CA:1483:A:H2'	1:CA:1484:C:O4'	2.20	0.41
1:CA:1493:A:H2'	1:CA:1494:G:OP2	2.20	0.41
1:CA:102:G:N3	1:CA:151:A:H2	2.18	0.41
1:CA:155:A:H2'	1:CA:156:C:O4'	2.19	0.41
1:CA:208:U:H2'	1:CA:210:C:C5	2.55	0.41
1:CA:392:C:H2'	1:CA:393:A:C8	2.54	0.41
1:CA:407:U:H3'	1:CA:408:A:H8	1.85	0.41
1:CA:839:C:O2'	1:CA:840:C:H5'	2.19	0.41
1:CA:949:A:H2'	1:CA:950:U:O4'	2.21	0.41
3:CD:151:GLN:HE21	3:CD:153:ARG:HB3	1.85	0.41
3:CD:171:GLU:O	3:CD:180:THR:HB	2.21	0.41
6:CG:65:LEU:C	6:CG:67:ASN:H	2.23	0.41
7:CH:28:SER:HB3	7:CH:58:LEU:HB2	2.01	0.41
8:CI:80:HIS:HE1	8:CI:103:VAL:O	2.03	0.41
10:CK:126:ARG:NH1	10:CK:126:ARG:HG2	2.35	0.41
10:CK:12:ARG:HB3	10:CK:13:LYS:HZ3	1.84	0.41
10:CK:22:ILE:HD13	10:CK:95:THR:CG2	2.49	0.41
10:CK:70:ALA:CA	10:CK:74:LYS:HB2	2.50	0.41
1:CA:562:U:H1'	11:CL:11:ARG:HB3	2.02	0.41
12:CM:18:LEU:HD23	12:CM:18:LEU:O	2.20	0.41
13:CN:21:ALA:O	13:CN:24:ALA:HB3	2.20	0.41
13:CN:82:LYS:HA	13:CN:82:LYS:HE2	2.01	0.41
13:CN:65:GLN:HB3	13:CN:82:LYS:HG2	2.01	0.41
22:DA:112:G:H2'	22:DA:113:C:C6	2.54	0.41
22:DA:30:C:H2'	22:DA:31:C:C5'	2.41	0.41
23:DB:1987:A:H2'	23:DB:1988:G:C8	2.54	0.41
23:DB:2016:U:H2'	23:DB:2017:U:C6	2.56	0.41
23:DB:212:G:H2'	23:DB:213:A:C8	2.55	0.41
23:DB:2720:U:H2'	23:DB:2721:A:C8	2.55	0.41
23:DB:2768:U:H2'	23:DB:2769:U:O4'	2.20	0.41
23:DB:2874:C:H2'	23:DB:2875:C:H6	1.84	0.41
23:DB:2886:A:C4	47:D0:27:LEU:HG	2.56	0.41
23:DB:340:A:H2'	23:DB:341:C:C5'	2.50	0.41
23:DB:910:A:N7	34:DM:16:ARG:HB3	2.35	0.41
23:DB:963:U:H2'	23:DB:964:C:H6	1.84	0.41
25:DC:27:LYS:N	25:DC:27:LYS:HD2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:42:ARG:CZ	25:DC:44:ASN:HB2	2.50	0.41
26:DD:106:LYS:HG3	26:DD:208:LYS:CE	2.51	0.41
28:DF:174:PHE:N	28:DF:175:PRO:CD	2.83	0.41
28:DF:38:GLY:N	28:DF:86:CYS:SG	2.93	0.41
28:DF:36:ASN:HD22	28:DF:87:LYS:H	1.68	0.41
28:DF:9:ASP:O	28:DF:10:GLU:CB	2.68	0.41
29:DG:96:ALA:HA	29:DG:124:CYS:SG	2.60	0.41
31:DJ:124:VAL:CG2	31:DJ:125:TYR:N	2.76	0.41
31:DJ:64:VAL:HG21	31:DJ:69:ARG:HG2	2.02	0.41
31:DJ:80:HIS:N	31:DJ:80:HIS:ND1	2.68	0.41
32:DK:17:ARG:HB2	32:DK:45:GLU:HB3	2.02	0.41
33:DL:100:ILE:O	33:DL:101:ILE:HD13	2.19	0.41
33:DL:116:VAL:O	33:DL:116:VAL:HG23	2.20	0.41
33:DL:79:LEU:N	33:DL:113:ALA:HB3	2.35	0.41
34:DM:15:GLY:O	34:DM:16:ARG:O	2.38	0.41
37:DP:40:GLN:O	37:DP:41:ALA:O	2.38	0.41
39:DR:18:GLN:H	39:DR:18:GLN:CD	2.21	0.41
39:DR:6:GLN:CB	39:DR:41:ILE:HD13	2.49	0.41
40:DS:7:HIS:ND1	40:DS:7:HIS:O	2.52	0.41
42:DU:71:ILE:CD1	42:DU:72:PHE:N	2.80	0.41
1:AA:112:G:O2'	1:AA:113:G:H5'	2.19	0.41
1:AA:1281:C:H3'	1:AA:1282:C:H6	1.85	0.41
1:AA:814:A:C5'	1:AA:1511:G:H4'	2.45	0.41
1:AA:243:A:C1'	1:AA:244:U:OP2	2.68	0.41
1:AA:254:G:H4'	16:AQ:19:SER:OG	2.20	0.41
1:AA:546:A:H4'	1:AA:548:G:O3'	2.21	0.41
1:AA:599:C:H2'	1:AA:600:A:H8	1.85	0.41
1:AA:728:A:H2'	1:AA:729:A:C8	2.56	0.41
1:AA:969:A:H2'	1:AA:970:C:O2	2.21	0.41
20:AB:107:ARG:HA	20:AB:110:ILE:CD1	2.50	0.41
20:AB:67:LEU:HD21	20:AB:157:PRO:CB	2.50	0.41
20:AB:185:ILE:HG22	20:AB:199:ILE:HB	2.02	0.41
3:AD:12:ARG:HG3	3:AD:33:ILE:HA	2.01	0.41
5:AF:42:TRP:CZ2	5:AF:61:LEU:HD23	2.55	0.41
6:AG:30:MET:HA	6:AG:38:ALA:CB	2.51	0.41
6:AG:49:LEU:HB3	6:AG:57:GLU:OE2	2.20	0.41
6:AG:8:GLN:HG2	6:AG:9:ARG:N	2.35	0.41
7:AH:113:ARG:NH2	7:AH:114:ALA:HA	2.21	0.41
8:AI:20:ILE:HD13	8:AI:85:ALA:CB	2.50	0.41
8:AI:56:MET:O	8:AI:58:GLU:N	2.53	0.41
9:AJ:96:VAL:O	9:AJ:98:VAL:HG13	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AL:21:PRO:HD2	11:AL:94:TYR:OH	2.20	0.41
12:AM:14:ALA:HB1	12:AM:33:LEU:HD11	2.02	0.41
17:AR:52:ARG:HH11	17:AR:52:ARG:CB	2.21	0.41
18:AS:78:THR:HG22	18:AS:79:TYR:N	2.35	0.41
50:B3:23:HIS:HB2	50:B3:24:LYS:H	1.66	0.41
50:B3:23:HIS:HE1	50:B3:51:LYS:HE2	1.84	0.41
23:BB:107:G:H2'	23:BB:108:G:H8	1.84	0.41
23:BB:1139:G:H5'	31:BJ:104:ALA:HB1	2.02	0.41
23:BB:118:A:H5'	23:BB:119:A:C8	2.41	0.41
23:BB:1601:G:H2'	23:BB:1602:U:O4'	2.20	0.41
23:BB:1949:G:H2'	23:BB:1950:G:C8	2.55	0.41
23:BB:2104:C:H4'	23:BB:2104:C:OP1	2.20	0.41
23:BB:2250:G:H21	23:BB:2496:C:H4'	1.85	0.41
23:BB:2491:U:H5''	23:BB:2570:G:H5'	2.00	0.41
23:BB:2665:A:H2'	23:BB:2666:C:O2	2.21	0.41
23:BB:2675:A:O2'	23:BB:2676:C:H5'	2.20	0.41
23:BB:2788:C:H2'	23:BB:2789:C:C6	2.55	0.41
23:BB:277:G:HO2'	23:BB:361:G:H1	1.68	0.41
23:BB:410:G:C2	23:BB:2407:A:C5	3.08	0.41
23:BB:583:G:H2'	23:BB:584:C:H6	1.85	0.41
23:BB:79:C:H2'	23:BB:80:G:H8	1.85	0.41
25:BC:222:THR:C	25:BC:224:MET:N	2.73	0.41
23:BB:1655:A:H5''	26:BD:123:LYS:HZ2	1.86	0.41
26:BD:133:THR:CG2	26:BD:134:HIS:N	2.83	0.41
27:BE:170:ARG:HH11	27:BE:170:ARG:HG2	1.85	0.41
27:BE:149:ILE:HG23	27:BE:183:PHE:CG	2.56	0.41
27:BE:48:THR:CG2	27:BE:49:ARG:N	2.80	0.41
28:BF:177:ARG:C	28:BF:177:ARG:NE	2.74	0.41
29:BG:129:GLU:C	29:BG:130:ILE:HD12	2.40	0.41
30:BH:44:ILE:O	30:BH:47:PHE:HD2	2.04	0.41
30:BH:53:GLU:C	30:BH:54:LEU:HD22	2.41	0.41
31:BJ:41:LYS:HD2	31:BJ:44:TYR:CB	2.47	0.41
32:BK:38:ILE:HA	32:BK:60:ALA:O	2.20	0.41
34:BM:108:VAL:C	34:BM:112:LEU:HD13	2.40	0.41
35:BN:24:MET:CE	35:BN:44:LEU:HD13	2.50	0.41
37:BP:92:ARG:N	37:BP:113:LEU:HD23	2.35	0.41
38:BQ:12:ARG:O	38:BQ:15:LYS:HB3	2.20	0.41
39:BR:40:MET:C	39:BR:54:VAL:HG11	2.41	0.41
41:BT:31:VAL:O	41:BT:83:ALA:O	2.38	0.41
41:BT:54:GLU:HB2	41:BT:91:GLN:HE22	1.85	0.41
23:BB:328:U:C4'	42:BU:65:GLN:HG3	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1162:C:H2'	1:CA:1163:A:H8	1.83	0.41
1:CA:1187:G:H3'	1:CA:1188:A:H8	1.86	0.41
1:CA:1206:G:H2'	1:CA:1207:G:O4'	2.20	0.41
1:CA:1469:C:H2'	1:CA:1470:U:O4'	2.19	0.41
1:CA:191:G:H2'	1:CA:192:A:H8	1.85	0.41
1:CA:343:U:O2'	1:CA:344:A:H2'	2.20	0.41
1:CA:42:G:O2'	1:CA:43:C:H5'	2.20	0.41
1:CA:489:C:H2'	1:CA:490:C:H6	1.86	0.41
1:CA:601:G:O2'	1:CA:602:A:H5'	2.21	0.41
20:CB:205:ALA:HB3	20:CB:208:ALA:CB	2.50	0.41
2:CC:22:PHE:CZ	9:CJ:11:LYS:HB3	2.55	0.41
3:CD:18:LEU:HD11	3:CD:59:LYS:HG3	2.02	0.41
4:CE:37:VAL:HA	4:CE:46:GLY:O	2.21	0.41
7:CH:4:ASP:OD1	7:CH:76:ARG:NH1	2.54	0.41
7:CH:72:GLU:O	7:CH:73:SER:HB2	2.20	0.41
8:CI:30:ASN:HD21	8:CI:66:VAL:N	2.18	0.41
9:CJ:32:THR:HG23	9:CJ:33:GLY:N	2.31	0.41
13:CN:56:PRO:HG2	13:CN:57:SER:H	1.85	0.41
9:CJ:66:GLU:O	13:CN:95:LEU:HA	2.20	0.41
17:CR:62:ARG:HB3	17:CR:69:TYR:CE1	2.56	0.41
18:CS:44:ILE:HG23	18:CS:62:THR:O	2.21	0.41
19:CT:43:LYS:HD3	19:CT:86:ALA:HB1	2.03	0.41
22:DA:55:U:H4'	28:DF:23:SER:OG	2.20	0.41
23:DB:1220:G:O2'	23:DB:1221:C:H5'	2.19	0.41
23:DB:123:G:H2'	23:DB:124:G:C8	2.55	0.41
23:DB:1306:C:H2'	23:DB:1307:A:H8	1.84	0.41
23:DB:1341:G:H2'	23:DB:1397:U:HO2'	1.85	0.41
23:DB:1399:C:H2'	23:DB:1400:U:H6	1.83	0.41
23:DB:1736:U:H2'	23:DB:1737:G:O4'	2.21	0.41
23:DB:1795:C:H2'	23:DB:1796:U:H6	1.85	0.41
23:DB:1869:G:HO2'	23:DB:1871:A:H62	1.62	0.41
23:DB:2251:G:OP2	23:DB:2251:G:H8	2.03	0.41
23:DB:2354:C:O2'	43:DW:31:LEU:HG	2.20	0.41
23:DB:2675:A:O2'	23:DB:2676:C:H5'	2.21	0.41
23:DB:2682:A:H1'	26:DD:22:ILE:HD13	2.02	0.41
23:DB:2788:C:H2'	23:DB:2789:C:C6	2.55	0.41
23:DB:532:A:H5''	38:DQ:27:ARG:NH1	2.34	0.41
23:DB:787:C:H3'	23:DB:791:C:H41	1.85	0.41
23:DB:941:A:H2'	23:DB:942:G:C8	2.55	0.41
25:DC:162:GLN:NE2	25:DC:174:ARG:NH2	2.68	0.41
23:DB:1789:A:H5'	25:DC:220:ARG:NH2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:DC:21:PRO:O	25:DC:22:GLU:CB	2.67	0.41
26:DD:106:LYS:O	26:DD:206:ALA:HB2	2.20	0.41
28:DF:100:GLU:C	28:DF:102:LEU:N	2.73	0.41
29:DG:12:ALA:C	29:DG:14:VAL:H	2.23	0.41
23:DB:2094:A:OP1	30:DH:22:LYS:HG3	2.21	0.41
31:DJ:4:PHE:HB2	31:DJ:5:THR:H	1.49	0.41
33:DL:135:ILE:N	33:DL:135:ILE:HD13	2.34	0.41
33:DL:38:GLN:HB2	33:DL:39:LYS:H	1.66	0.41
33:DL:38:GLN:HB2	33:DL:39:LYS:HZ3	1.85	0.41
35:DN:12:ARG:NH1	35:DN:16:HIS:HD2	2.18	0.41
35:DN:86:ARG:NH1	35:DN:94:TYR:OH	2.53	0.41
23:DB:533:G:C5'	38:DQ:23:TYR:CD1	3.02	0.41
40:DS:75:PHE:O	40:DS:103:ILE:O	2.37	0.41
40:DS:7:HIS:O	40:DS:8:ARG:HB2	2.20	0.41
41:DT:48:GLN:CG	41:DT:49:LYS:N	2.84	0.41
42:DU:27:VAL:CG1	42:DU:33:VAL:HG22	2.50	0.41
42:DU:69:VAL:CG1	42:DU:77:GLY:HA2	2.50	0.41
43:DW:74:LYS:HB3	43:DW:75:ASN:H	1.55	0.41
1:AA:1085:U:H3'	1:AA:1086:U:H5	1.86	0.41
1:AA:1179:A:O2'	1:AA:1180:A:H5'	2.20	0.41
1:AA:1263:C:H2'	1:AA:1264:U:C6	2.55	0.41
1:AA:1350:A:H2'	1:AA:1351:U:C6	2.55	0.41
1:AA:136:C:H1'	15:AP:1:MET:HE1	2.02	0.41
1:AA:1494:G:H2'	1:AA:1495:U:C6	2.54	0.41
1:AA:177:G:P	19:AT:59:ARG:HE	2.43	0.41
1:AA:255:G:H1'	16:AQ:17:GLU:OE1	2.20	0.41
1:AA:285:C:H2'	1:AA:286:C:C6	2.56	0.41
1:AA:313:A:H2'	1:AA:314:C:H6	1.82	0.41
1:AA:596:A:O2'	1:AA:597:G:H5'	2.20	0.41
1:AA:746:A:H2'	1:AA:747:A:H8	1.86	0.41
20:AB:162:VAL:HG13	20:AB:184:ALA:CB	2.47	0.41
2:AC:179:ALA:HB1	2:AC:202:PHE:CE1	2.40	0.41
4:AE:17:VAL:HB	4:AE:34:ALA:HB2	2.03	0.41
4:AE:45:VAL:HG23	4:AE:71:ILE:CG2	2.50	0.41
6:AG:142:ARG:HB2	6:AG:142:ARG:NH1	2.35	0.41
7:AH:87:ARG:O	7:AH:91:LEU:HG	2.21	0.41
8:AI:84:ARG:NH1	8:AI:84:ARG:HG2	2.34	0.41
9:AJ:48:ARG:HG2	9:AJ:48:ARG:NH1	2.35	0.41
11:AL:120:ARG:NH1	11:AL:120:ARG:HG3	2.36	0.41
14:AO:57:ARG:HH11	14:AO:57:ARG:HG3	1.85	0.41
15:AP:6:LEU:CD1	15:AP:71:VAL:HB	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AU:36:PHE:HB3	21:AU:40:PRO:HG3	2.00	0.41
49:B2:3:ARG:O	49:B2:5:PHE:N	2.53	0.41
51:B4:33:HIS:N	51:B4:33:HIS:CD2	2.79	0.41
23:BB:1107:G:O2'	23:BB:1108:U:H5'	2.20	0.41
23:BB:2386:A:H2'	23:BB:2387:U:H6	1.82	0.41
23:BB:2389:G:H5''	23:BB:2390:U:O4'	2.20	0.41
23:BB:2425:A:H4'	23:BB:2426:A:OP2	2.19	0.41
23:BB:2527:C:O2'	51:B4:1:MET:HA	2.20	0.41
23:BB:289:G:H2'	23:BB:290:U:O4'	2.21	0.41
25:BC:209:ALA:C	25:BC:213:ARG:HD2	2.41	0.41
23:BB:1797:G:P	25:BC:251:THR:HG23	2.60	0.41
25:BC:52:HIS:CD2	25:BC:53:ILE:HG13	2.56	0.41
26:BD:124:ARG:NH2	26:BD:163:GLY:HA3	2.35	0.41
26:BD:159:LYS:HD3	26:BD:159:LYS:C	2.41	0.41
26:BD:118:PHE:CA	26:BD:164:GLN:HG3	2.50	0.41
26:BD:166:GLY:O	26:BD:167:ASN:C	2.58	0.41
27:BE:6:LYS:NZ	27:BE:119:ILE:H	2.18	0.41
27:BE:141:MET:HB2	27:BE:143:LEU:HD22	2.02	0.41
27:BE:48:THR:CG2	27:BE:86:ALA:HB1	2.50	0.41
28:BF:56:LEU:HD23	28:BF:56:LEU:O	2.20	0.41
31:BJ:80:HIS:O	31:BJ:81:ILE:C	2.59	0.41
23:BB:1131:G:C1'	31:BJ:85:LYS:HZ2	2.34	0.41
32:BK:119:ALA:O	32:BK:121:GLU:N	2.53	0.41
32:BK:34:GLY:O	32:BK:36:GLY:N	2.53	0.41
32:BK:87:LEU:CA	32:BK:95:ILE:HG22	2.50	0.41
33:BL:78:ARG:HH21	33:BL:82:LEU:HB2	1.83	0.41
34:BM:13:HIS:O	34:BM:14:LYS:HB3	2.21	0.41
34:BM:7:THR:CG2	34:BM:8:LYS:N	2.82	0.41
35:BN:13:ASN:C	35:BN:15:SER:N	2.74	0.41
36:BO:45:SER:C	36:BO:47:VAL:H	2.23	0.41
23:BB:2693:G:H21	37:BP:95:LYS:NZ	2.18	0.41
38:BQ:59:LEU:O	38:BQ:62:ALA:HB3	2.19	0.41
38:BQ:60:TRP:CE2	38:BQ:93:ILE:HA	2.54	0.41
39:BR:31:GLU:HB3	39:BR:32:THR:H	1.53	0.41
39:BR:79:ARG:HD3	39:BR:86:GLN:CD	2.40	0.41
40:BS:12:SER:O	40:BS:16:LYS:HB2	2.21	0.41
46:BZ:3:LYS:O	46:BZ:3:LYS:HG3	2.19	0.41
1:CA:107:G:C2'	1:CA:108:G:H5''	2.50	0.41
1:CA:1184:G:C2	1:CA:1185:G:C8	3.07	0.41
1:CA:1413:A:O2'	1:CA:1414:U:H5'	2.20	0.41
1:CA:1434:A:H2'	1:CA:1435:G:O4'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:285:C:H2'	1:CA:286:C:C6	2.55	0.41
1:CA:580:C:H2'	1:CA:581:G:O4'	2.21	0.41
1:CA:605:U:H2'	1:CA:606:G:H8	1.86	0.41
1:CA:664:G:N2	1:CA:741:G:H1	2.03	0.41
1:CA:813:U:O2'	1:CA:814:A:H5'	2.20	0.41
1:CA:868:C:H2'	1:CA:869:G:O4'	2.20	0.41
20:CB:125:PHE:C	20:CB:127:LYS:H	2.23	0.41
2:CC:143:LEU:HB2	2:CC:144:GLY:H	1.67	0.41
3:CD:157:ALA:HA	3:CD:160:LEU:CD1	2.47	0.41
3:CD:155:LYS:C	3:CD:157:ALA:H	2.23	0.41
3:CD:160:LEU:HD22	3:CD:160:LEU:C	2.41	0.41
3:CD:195:ASN:HB3	3:CD:197:HIS:NE2	2.35	0.41
3:CD:27:ILE:O	3:CD:27:ILE:HG22	2.20	0.41
3:CD:74:TYR:CD1	3:CD:92:LEU:HD12	2.55	0.41
6:CG:16:LYS:HB3	6:CG:17:PHE:CD1	2.54	0.41
1:CA:1180:A:OP1	8:CI:104:THR:HG23	2.21	0.41
8:CI:98:ARG:HA	8:CI:103:VAL:CG2	2.45	0.41
9:CJ:83:THR:HG23	9:CJ:87:LEU:HD23	2.01	0.41
13:CN:50:LEU:N	13:CN:51:PRO:CD	2.68	0.41
1:CA:1220:G:P	13:CN:52:ARG:HH22	2.43	0.41
23:DB:2884:U:N3	47:D0:48:TYR:CE1	2.88	0.41
49:D2:17:GLY:C	49:D2:18:PHE:CG	2.93	0.41
23:DB:1099:G:P	52:DI:3:LYS:CA	2.95	0.41
23:DB:1372:U:H2'	23:DB:1373:A:H8	1.85	0.41
23:DB:1601:G:H2'	23:DB:1602:U:O4'	2.20	0.41
23:DB:1612:C:O3'	49:D2:5:PHE:HD2	2.02	0.41
23:DB:1685:C:H2'	23:DB:1686:C:C6	2.55	0.41
23:DB:1637:A:H5'	23:DB:1760:C:O2'	2.20	0.41
23:DB:1949:G:H2'	23:DB:1950:G:C8	2.55	0.41
23:DB:1998:A:H2'	23:DB:1999:C:C6	2.55	0.41
23:DB:2026:U:C2	23:DB:2027:G:C8	3.07	0.41
23:DB:2321:U:H5''	23:DB:2321:U:O2	2.19	0.41
23:DB:2515:C:OP1	31:DJ:82:GLY:C	2.59	0.41
23:DB:2641:G:H5''	31:DJ:78:THR:HB	2.01	0.41
23:DB:2849:U:N3	23:DB:2867:G:H1'	2.35	0.41
23:DB:491:G:C2	23:DB:492:A:H1'	2.55	0.41
23:DB:853:C:H2'	23:DB:854:C:C6	2.54	0.41
25:DC:10:PRO:CD	25:DC:202:ARG:NH2	2.80	0.41
25:DC:143:VAL:HG12	25:DC:144:GLU:N	2.27	0.41
25:DC:224:MET:HB3	25:DC:225:ASN:H	1.70	0.41
23:DB:2228:G:OP1	25:DC:257:ARG:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:DD:41:ALA:C	26:DD:43:ASP:H	2.23	0.41
26:DD:28:GLU:CD	26:DD:97:SER:HA	2.41	0.41
27:DE:108:ILE:HG13	27:DE:109:LEU:N	2.29	0.41
28:DF:34:THR:O	28:DF:153:ILE:HA	2.20	0.41
28:DF:63:LYS:HD2	28:DF:64:PRO:HD2	2.02	0.41
29:DG:36:LEU:HD22	29:DG:40:VAL:CG1	2.50	0.41
29:DG:95:ALA:HB1	29:DG:130:ILE:CD1	2.48	0.41
30:DH:57:LYS:H	30:DH:57:LYS:HG3	1.60	0.41
33:DL:104:GLN:HB2	33:DL:105:ILE:H	1.60	0.41
33:DL:90:VAL:HG22	33:DL:90:VAL:O	2.20	0.41
35:DN:36:THR:HG22	35:DN:40:LYS:HD2	2.02	0.41
35:DN:49:GLU:N	35:DN:50:PRO:HD2	2.35	0.41
36:DO:25:ARG:HB2	36:DO:40:ILE:HG12	2.03	0.41
37:DP:91:VAL:CG2	37:DP:92:ARG:H	2.25	0.41
38:DQ:25:GLY:HA2	38:DQ:29:ARG:CD	2.48	0.41
38:DQ:89:ILE:H	38:DQ:89:ILE:CD1	2.30	0.41
39:DR:47:VAL:HG13	39:DR:48:LYS:N	2.34	0.41
40:DS:72:THR:CG2	40:DS:106:VAL:HG23	2.50	0.41
43:DW:21:GLY:HA2	43:DW:25:PHE:CD1	2.54	0.41
43:DW:42:THR:CB	43:DW:75:ASN:ND2	2.83	0.41
45:DY:43:ILE:HA	45:DY:46:MET:HB2	2.01	0.41
46:DZ:55:GLY:HA2	46:DZ:59:ARG:CB	2.31	0.41
46:DZ:66:ILE:HB	46:DZ:67:PRO:CD	2.42	0.41
1:AA:1110:A:H2'	1:AA:1111:A:C8	2.49	0.41
1:AA:1157:A:H4'	1:AA:1158:C:O5'	2.20	0.41
1:AA:119:A:H4'	1:AA:120:A:O4'	2.21	0.41
1:AA:132:C:H5''	19:AT:68:LYS:HD2	2.02	0.41
1:AA:1340:A:O2'	1:AA:1341:U:H5'	2.20	0.41
1:AA:224:U:H2'	1:AA:225:C:C6	2.56	0.41
1:AA:605:U:H2'	1:AA:606:G:H8	1.86	0.41
1:AA:635:A:H2'	1:AA:636:U:H6	1.85	0.41
1:AA:813:U:O2'	1:AA:814:A:H5'	2.20	0.41
1:AA:830:G:H2'	1:AA:831:A:H8	1.85	0.41
1:AA:903:G:H2'	1:AA:904:U:C6	2.55	0.41
1:AA:75:G:N1	1:AA:96:U:C4	2.88	0.41
2:AC:18:ASN:O	2:AC:55:VAL:HA	2.20	0.41
3:AD:123:MET:HG3	3:AD:127:ARG:C	2.40	0.41
3:AD:163:GLN:HB2	3:AD:164:ARG:HH12	1.82	0.41
6:AG:142:ARG:C	6:AG:146:ALA:HB3	2.41	0.41
1:AA:600:A:H5''	7:AH:88:LYS:HD2	2.02	0.41
11:AL:79:ILE:HD13	11:AL:96:THR:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AP:12:LYS:C	15:AP:14:ARG:H	2.24	0.41
18:AS:64:GLU:OE1	18:AS:65:MET:HG3	2.20	0.41
19:AT:19:HIS:CE1	19:AT:23:ARG:HG3	2.55	0.41
21:AU:40:PRO:HG2	21:AU:41:THR:N	2.28	0.41
51:B4:15:LYS:N	51:B4:15:LYS:HD2	2.35	0.41
23:BB:1131:G:C1'	23:BB:1133:A:H62	2.24	0.41
23:BB:1297:C:OP1	23:BB:2710:C:H4'	2.20	0.41
23:BB:175:G:H2'	23:BB:176:A:H8	1.83	0.41
23:BB:1841:U:H2'	23:BB:1842:G:C8	2.55	0.41
23:BB:1924:C:O2'	23:BB:1925:C:H5'	2.21	0.41
23:BB:1998:A:H2'	23:BB:1999:C:H6	1.84	0.41
23:BB:2034:U:O2'	23:BB:2035:G:H5'	2.21	0.41
23:BB:1655:A:H2	23:BB:2049:G:O3'	2.03	0.41
23:BB:2185:U:H2'	23:BB:2186:G:C8	2.51	0.41
23:BB:2341:G:H2'	23:BB:2342:C:C6	2.55	0.41
23:BB:2469:A:C5'	34:BM:55:ARG:HG3	2.50	0.41
23:BB:2649:C:O2'	23:BB:2650:U:H5'	2.19	0.41
23:BB:2677:G:H2'	23:BB:2678:C:C6	2.55	0.41
23:BB:2782:G:O2'	23:BB:2783:U:H5'	2.20	0.41
23:BB:299:A:H2	23:BB:319:G:N3	2.18	0.41
23:BB:482:A:N6	23:BB:506:G:H1'	2.35	0.41
23:BB:527:C:C5'	23:BB:2779:U:H3	2.34	0.41
23:BB:730:A:H3'	55:BB:3597:HOH:O	2.19	0.41
23:BB:795:C:H2'	23:BB:796:C:C6	2.55	0.41
23:BB:962:G:H2'	23:BB:963:U:H6	1.85	0.41
25:BC:156:SER:O	25:BC:194:VAL:O	2.39	0.41
25:BC:212:TRP:C	25:BC:214:GLY:H	2.24	0.41
23:BB:778:G:C5'	25:BC:48:ILE:HD12	2.51	0.41
26:BD:24:VAL:CG2	26:BD:193:VAL:HG11	2.27	0.41
27:BE:114:ARG:HG2	27:BE:114:ARG:H	1.65	0.41
27:BE:116:ASP:HB2	27:BE:185:LYS:CA	2.36	0.41
27:BE:52:VAL:C	27:BE:54:GLY:N	2.73	0.41
27:BE:84:THR:OG1	27:BE:85:PHE:N	2.53	0.41
28:BF:52:ALA:HA	28:BF:148:VAL:CG1	2.50	0.41
28:BF:34:THR:CG2	28:BF:89:THR:HG22	2.48	0.41
52:BI:2:LYS:NZ	52:BI:2:LYS:HB3	2.36	0.41
31:BJ:102:GLU:HG3	31:BJ:124:VAL:HG21	2.02	0.41
32:BK:33:ALA:CB	32:BK:39:ILE:HD11	2.50	0.41
33:BL:60:ARG:HD3	50:B3:11:LYS:HD2	2.02	0.41
33:BL:6:LEU:HD22	33:BL:6:LEU:N	2.35	0.41
35:BN:9:GLN:HE21	35:BN:17:ARG:CZ	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BN:22:ARG:HE	35:BN:69:ARG:HE	1.67	0.41
36:BO:76:LYS:HA	36:BO:76:LYS:HD2	1.95	0.41
38:BQ:21:LYS:C	38:BQ:23:TYR:H	2.23	0.41
38:BQ:23:TYR:O	38:BQ:24:TYR:HB2	2.20	0.41
41:BT:12:ARG:HE	44:BX:29:ARG:NH2	2.11	0.41
24:BV:51:GLN:NE2	24:BV:57:TYR:OH	2.53	0.41
45:BY:3:THR:O	45:BY:6:ILE:HD12	2.20	0.41
46:BZ:58:ASP:HB3	46:BZ:63:ARG:HG2	2.01	0.41
1:CA:1032:G:H2'	1:CA:1033:G:O4'	2.20	0.41
1:CA:1190:G:H5'	2:CC:175:HIS:HE2	1.84	0.41
1:CA:1438:G:O2'	1:CA:1439:G:H5'	2.20	0.41
1:CA:224:U:H2'	1:CA:225:C:C6	2.55	0.41
1:CA:340:U:H2'	1:CA:341:C:C6	2.53	0.41
1:CA:97:G:H3'	1:CA:98:A:H8	1.85	0.41
20:CB:101:THR:HG23	20:CB:102:ASN:N	2.36	0.41
20:CB:72:LYS:HD2	20:CB:164:ASP:HB3	2.02	0.41
2:CC:4:VAL:HG22	2:CC:5:HIS:N	2.35	0.41
3:CD:32:LYS:HG3	3:CD:32:LYS:O	2.20	0.41
3:CD:60:VAL:HA	3:CD:63:ILE:CG2	2.51	0.41
6:CG:99:ALA:O	6:CG:103:ILE:HG13	2.20	0.41
7:CH:100:ILE:HD12	7:CH:100:ILE:C	2.41	0.41
13:CN:51:PRO:CG	13:CN:52:ARG:H	2.30	0.41
13:CN:89:ARG:CB	13:CN:89:ARG:HH11	2.34	0.41
14:CO:80:LEU:CD2	14:CO:84:LEU:HD22	2.49	0.41
15:CP:44:SER:C	15:CP:46:LYS:N	2.73	0.41
19:CT:66:ILE:HG23	19:CT:70:LYS:CG	2.51	0.41
49:D2:7:PRO:O	49:D2:8:SER:CB	2.67	0.41
50:D3:35:LYS:HG3	50:D3:35:LYS:H	1.70	0.41
23:DB:247:G:C6	50:D3:7:ARG:HD2	2.56	0.41
22:DA:52:A:H2'	22:DA:53:A:C8	2.53	0.41
23:DB:1023:U:H2'	23:DB:1024:G:H5'	2.02	0.41
23:DB:1205:A:H1'	23:DB:1206:G:P	2.61	0.41
23:DB:1281:G:H2'	23:DB:1282:U:C6	2.56	0.41
23:DB:2073:C:O2'	23:DB:2074:U:H5'	2.20	0.41
23:DB:2090:A:C2'	46:DZ:49:ARG:NE	2.83	0.41
23:DB:2184:A:H2'	23:DB:2185:U:C6	2.56	0.41
23:DB:2273:A:H2'	23:DB:2274:A:C8	2.56	0.41
23:DB:2511:U:O5'	23:DB:2511:U:H6	2.02	0.41
23:DB:264:C:C2'	23:DB:265:A:H5''	2.50	0.41
23:DB:2659:G:N2	23:DB:2661:G:H5''	2.36	0.41
23:DB:2677:G:H2'	23:DB:2678:C:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:2732:G:H3'	23:DB:2733:A:H5'	2.02	0.41
23:DB:300:A:H2'	23:DB:334:C:H1'	2.01	0.41
23:DB:365:U:H6	23:DB:365:U:O5'	2.02	0.41
23:DB:484:C:H2'	23:DB:485:C:H6	1.84	0.41
23:DB:857:G:H2'	23:DB:858:G:C5'	2.48	0.41
25:DC:252:LYS:HB3	25:DC:253:GLY:H	1.69	0.41
25:DC:87:SER:N	25:DC:155:ARG:NH1	2.68	0.41
26:DD:10:GLY:HA2	26:DD:23:PRO:O	2.21	0.41
26:DD:34:VAL:H	26:DD:90:PHE:CA	2.33	0.41
28:DF:34:THR:HB	28:DF:154:THR:O	2.21	0.41
29:DG:132:LEU:O	29:DG:132:LEU:HD12	2.20	0.41
52:DI:100:ILE:O	52:DI:139:VAL:HG13	2.20	0.41
31:DJ:61:LYS:O	31:DJ:62:VAL:HG13	2.21	0.41
31:DJ:76:HIS:HB2	31:DJ:86:GLN:CD	2.41	0.41
32:DK:19:VAL:O	32:DK:41:ILE:HD11	2.20	0.41
34:DM:35:ALA:HB2	34:DM:100:LYS:NZ	2.35	0.41
35:DN:11:ASN:HA	35:DN:11:ASN:HD22	1.70	0.41
36:DO:55:GLU:O	36:DO:56:LYS:C	2.57	0.41
37:DP:38:ARG:HG3	37:DP:39:LEU:N	2.20	0.41
23:DB:1187:G:H5''	39:DR:86:GLN:HE22	1.85	0.41
41:DT:66:LYS:HA	41:DT:76:ARG:HE	1.85	0.41
41:DT:84:TYR:O	41:DT:85:VAL:HG22	2.20	0.41
22:DA:82:U:H5''	45:DY:16:LEU:HD11	2.03	0.41
1:AA:1005:A:O2'	1:AA:1006:G:H5'	2.21	0.41
1:AA:1347:G:N2	1:AA:1373:G:H2'	2.35	0.41
1:AA:1370:G:O2'	1:AA:1371:G:H5'	2.20	0.41
1:AA:1395:C:H2'	1:AA:1395:C:O2	2.21	0.41
1:AA:265:G:H4'	16:AQ:66:LEU:O	2.20	0.41
1:AA:404:G:H2'	1:AA:405:U:H6	1.85	0.41
1:AA:489:C:H2'	1:AA:490:C:H6	1.86	0.41
1:AA:551:U:H2'	1:AA:552:U:H6	1.86	0.41
1:AA:666:G:H5'	1:AA:726:C:H1'	2.03	0.41
1:AA:66:A:O2'	1:AA:67:C:H5'	2.21	0.41
1:AA:68:G:H5'	1:AA:171:A:O2'	2.21	0.41
1:AA:738:C:H2'	1:AA:739:C:H6	1.86	0.41
20:AB:45:THR:HA	20:AB:48:MET:CG	2.51	0.41
2:AC:14:VAL:HG23	2:AC:15:LYS:HG2	2.03	0.41
2:AC:7:ASN:OD1	2:AC:183:TYR:HB3	2.21	0.41
3:AD:32:LYS:HB3	3:AD:35:GLN:CG	2.50	0.41
4:AE:95:MET:CE	4:AE:114:LEU:HD11	2.51	0.41
6:AG:21:LEU:HG	6:AG:22:LEU:H	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:22:LEU:O	6:AG:26:VAL:HG13	2.19	0.41
10:AK:122:PRO:CG	21:AU:35:GLU:HG2	2.51	0.41
13:AN:20:PHE:CG	13:AN:24:ALA:HB2	2.56	0.41
18:AS:48:ILE:HG22	18:AS:49:ALA:N	2.35	0.41
21:AU:33:ARG:CZ	21:AU:34:ARG:HG3	2.50	0.41
21:AU:33:ARG:HB3	21:AU:34:ARG:H	1.64	0.41
47:B0:32:THR:HG21	47:B0:35:GLU:HB3	2.02	0.41
49:B2:21:ARG:CB	49:B2:31:LEU:HD21	2.49	0.41
51:B4:7:VAL:HB	51:B4:35:GLN:HE21	1.85	0.41
23:BB:1281:G:H2'	23:BB:1282:U:C6	2.56	0.41
23:BB:1299:G:C5'	23:BB:1301:A:H1'	2.51	0.41
23:BB:1498:C:H2'	23:BB:1499:C:H6	1.86	0.41
23:BB:15:G:H2'	23:BB:16:C:H6	1.84	0.41
23:BB:2385:C:H2'	23:BB:2386:A:H8	1.86	0.41
23:BB:2457:U:C2'	23:BB:2458:G:H5'	2.49	0.41
23:BB:2710:C:H2'	23:BB:2711:A:C8	2.56	0.41
23:BB:2825:G:C2'	23:BB:2826:A:H5'	2.51	0.41
23:BB:477:A:H2'	23:BB:478:A:O4'	2.20	0.41
23:BB:724:U:O2'	23:BB:725:G:H5'	2.21	0.41
23:BB:466:A:H2	23:BB:795:C:O2	2.03	0.41
23:BB:990:A:N3	39:BR:79:ARG:NH2	2.68	0.41
25:BC:265:PHE:HB3	25:BC:266:ILE:H	1.74	0.41
26:BD:109:VAL:O	26:BD:172:VAL:HG22	2.21	0.41
26:BD:89:GLU:C	26:BD:91:THR:H	2.23	0.41
28:BF:133:GLU:CD	28:BF:150:GLY:H	2.24	0.41
28:BF:133:GLU:OE1	28:BF:150:GLY:N	2.53	0.41
28:BF:87:LYS:O	28:BF:88:VAL:HB	2.20	0.41
29:BG:29:ASN:N	29:BG:29:ASN:ND2	2.68	0.41
23:BB:2198:A:N6	30:BH:25:TYR:HE1	2.18	0.41
30:BH:31:VAL:HG22	30:BH:37:VAL:HG22	2.02	0.41
31:BJ:109:LEU:O	31:BJ:111:LYS:HG3	2.21	0.41
23:BB:1131:G:C1'	31:BJ:85:LYS:NZ	2.83	0.41
32:BK:20:MET:HB3	32:BK:42:THR:HB	2.01	0.41
27:BE:25:GLU:CD	33:BL:17:LYS:HG2	2.40	0.41
33:BL:21:ARG:CZ	33:BL:21:ARG:HB3	2.50	0.41
33:BL:60:ARG:O	33:BL:60:ARG:HG3	2.20	0.41
39:BR:42:ALA:C	39:BR:44:GLY:H	2.23	0.41
39:BR:78:ARG:HD3	39:BR:87:GLN:C	2.40	0.41
40:BS:51:LEU:HD13	40:BS:51:LEU:C	2.41	0.41
41:BT:70:HIS:CG	41:BT:74:ILE:HB	2.55	0.41
45:BY:24:LEU:HD23	45:BY:29:ARG:HH22	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1133:G:H2'	1:CA:1134:G:O4'	2.21	0.41
1:CA:1363:A:O2'	1:CA:1365:G:N7	2.45	0.41
1:CA:371:A:H2'	1:CA:372:C:C6	2.54	0.41
1:CA:644:U:O2'	1:CA:645:G:H5'	2.20	0.41
1:CA:79:G:H2'	1:CA:80:A:H8	1.83	0.41
1:CA:828:U:O2'	20:CB:24:PRO:HB3	2.20	0.41
3:CD:158:LEU:CD1	3:CD:177:MET:HG2	2.50	0.41
3:CD:93:LEU:HA	3:CD:96:ARG:HG3	2.03	0.41
8:CI:15:ALA:O	8:CI:66:VAL:HA	2.21	0.41
9:CJ:57:VAL:HG22	9:CJ:58:ASN:N	2.26	0.41
9:CJ:73:LEU:C	9:CJ:73:LEU:HD13	2.41	0.41
10:CK:82:GLU:HB3	10:CK:108:ASN:HB3	2.02	0.41
11:CL:29:LYS:O	11:CL:81:ILE:N	2.50	0.41
14:CO:30:LEU:O	14:CO:34:GLN:HG3	2.20	0.41
23:DB:1345:C:H5'	23:DB:1396:U:H5	1.85	0.41
23:DB:1394:U:O2'	23:DB:1395:A:H5'	2.21	0.41
23:DB:1434:A:OP1	23:DB:1434:A:H4'	2.20	0.41
23:DB:179:C:OP1	49:D2:28:ARG:HD3	2.18	0.41
23:DB:1977:A:H2'	23:DB:1978:A:O4'	2.21	0.41
23:DB:2034:U:O2'	23:DB:2035:G:H5'	2.20	0.41
23:DB:2070:A:H2'	23:DB:2071:A:O4'	2.20	0.41
23:DB:2072:C:H2'	23:DB:2073:C:C6	2.54	0.41
23:DB:2101:A:O2'	23:DB:2102:G:H5'	2.21	0.41
23:DB:2092:U:H5	23:DB:2226:C:OP2	2.02	0.41
23:DB:2642:G:H5''	31:DJ:80:HIS:HB2	2.02	0.41
23:DB:281:C:H2'	23:DB:282:A:C8	2.56	0.41
23:DB:2849:U:O4'	23:DB:2868:A:C6	2.73	0.41
23:DB:374:A:H2'	23:DB:375:G:O4'	2.21	0.41
23:DB:866:A:N1	23:DB:913:U:H4'	2.36	0.41
23:DB:928:A:H1'	45:DY:1:ALA:CA	2.48	0.41
23:DB:2204:G:C5'	25:DC:149:LYS:HE3	2.50	0.41
25:DC:160:TYR:CE2	25:DC:193:GLU:HG2	2.56	0.41
23:DB:1903:G:OP1	25:DC:239:PHE:CD2	2.74	0.41
23:DB:2032:G:N2	26:DD:150:GLN:HB3	2.36	0.41
27:DE:135:ALA:O	27:DE:139:LYS:N	2.53	0.41
28:DF:136:ILE:HG22	28:DF:137:PHE:N	2.35	0.41
29:DG:104:LEU:HD12	29:DG:106:LEU:HD21	2.02	0.41
29:DG:132:LEU:C	29:DG:132:LEU:HD12	2.40	0.41
30:DH:127:GLU:HB2	30:DH:143:ILE:CG2	2.50	0.41
23:DB:6:A:H4'	31:DJ:133:ALA:C	2.41	0.41
31:DJ:38:GLY:O	31:DJ:41:LYS:HD3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DJ:48:VAL:HG12	31:DJ:49:ASP:N	2.35	0.41
33:DL:143:GLU:HG2	33:DL:143:GLU:O	2.20	0.41
34:DM:129:THR:O	34:DM:131:VAL:HG12	2.20	0.41
34:DM:71:LYS:HZ3	34:DM:71:LYS:CA	2.29	0.41
35:DN:72:ASP:OD2	35:DN:75:ILE:HG23	2.20	0.41
35:DN:75:ILE:HG13	35:DN:76:VAL:N	2.36	0.41
36:DO:26:LEU:HD22	36:DO:93:ASP:HA	2.02	0.41
37:DP:50:ARG:CB	37:DP:50:ARG:HH11	2.31	0.41
39:DR:18:GLN:O	39:DR:19:THR:HG23	2.19	0.41
39:DR:27:ILE:HG23	39:DR:27:ILE:O	2.20	0.41
42:DU:42:LYS:N	42:DU:57:ILE:HD12	2.28	0.41
43:DW:3:LYS:HA	43:DW:3:LYS:HZ2	1.79	0.41
43:DW:57:THR:O	43:DW:58:LEU:C	2.59	0.41
45:DY:2:LYS:H	45:DY:37:ARG:HB2	1.83	0.41
46:DZ:28:VAL:HG23	46:DZ:29:GLY:H	1.85	0.41
46:DZ:34:LEU:HB3	46:DZ:35:ASP:H	1.54	0.41
1:AA:1217:C:H2'	1:AA:1218:C:C6	2.55	0.41
1:AA:1300:G:C2'	1:AA:1301:U:OP2	2.69	0.41
1:AA:1457:G:H2'	1:AA:1458:G:H8	1.86	0.41
1:AA:407:U:H3'	1:AA:408:A:H8	1.85	0.41
1:AA:878:A:H2'	1:AA:879:C:C6	2.56	0.41
1:AA:91:U:C4	1:AA:92:U:C4	3.09	0.41
1:AA:92:U:O2'	1:AA:93:U:H5'	2.20	0.41
1:AA:989:U:H2'	1:AA:990:C:H6	1.85	0.41
20:AB:76:SER:OG	20:AB:92:ASN:HB2	2.21	0.41
2:AC:35:ASP:O	2:AC:38:VAL:HG22	2.21	0.41
2:AC:86:LEU:O	2:AC:90:VAL:HG22	2.20	0.41
3:AD:109:THR:HG23	3:AD:112:GLU:H	1.86	0.41
3:AD:176:LYS:O	3:AD:177:MET:C	2.59	0.41
6:AG:10:LYS:NZ	6:AG:10:LYS:HB2	2.36	0.41
8:AI:24:ASN:C	8:AI:26:LYS:HZ2	2.24	0.41
10:AK:95:THR:HG23	10:AK:96:ILE:HG13	2.01	0.41
14:AO:58:MET:HG2	14:AO:58:MET:H	1.68	0.41
15:AP:40:ASN:HD21	15:AP:42:ILE:CG1	2.33	0.41
15:AP:44:SER:O	15:AP:45:GLU:C	2.59	0.41
18:AS:40:PHE:CG	18:AS:41:PRO:HD2	2.56	0.41
19:AT:54:GLN:N	19:AT:55:PRO:CD	2.82	0.41
23:BB:2372:U:H1'	48:B1:45:HIS:CE1	2.56	0.41
50:B3:34:LYS:HB3	50:B3:35:LYS:CE	2.50	0.41
51:B4:10:LEU:O	51:B4:11:CYS:HB3	2.21	0.41
22:BA:12:C:H41	43:BW:50:VAL:HG23	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:BA:81:G:H2'	22:BA:82:U:C6	2.55	0.41
23:BB:1097:U:H2'	23:BB:1098:A:C5'	2.50	0.41
23:BB:1036:G:C6	23:BB:1120:G:C6	3.07	0.41
23:BB:32:C:H5'	23:BB:1238:G:OP1	2.19	0.41
23:BB:1372:U:H2'	23:BB:1373:A:C8	2.55	0.41
23:BB:1593:A:H2'	23:BB:1594:U:O4'	2.20	0.41
23:BB:1945:G:H2'	23:BB:1946:U:H6	1.81	0.41
23:BB:1133:A:N6	23:BB:2025:C:O2'	2.54	0.41
23:BB:2205:A:H2'	23:BB:2206:C:H6	1.86	0.41
23:BB:2659:G:N2	23:BB:2661:G:H5''	2.36	0.41
23:BB:2668:G:O2'	23:BB:2669:G:H5'	2.20	0.41
23:BB:484:C:H2'	23:BB:485:C:H6	1.85	0.41
23:BB:490:C:H4'	23:BB:491:G:OP2	2.20	0.41
23:BB:545:U:C2	23:BB:548:G:OP2	2.74	0.41
23:BB:599:A:H2'	23:BB:600:G:H8	1.84	0.41
26:BD:138:LEU:CD2	26:BD:138:LEU:N	2.79	0.41
26:BD:138:LEU:HD23	26:BD:138:LEU:N	2.26	0.41
26:BD:141:ARG:O	26:BD:142:VAL:C	2.58	0.41
26:BD:165:MET:CE	26:BD:166:GLY:H	2.31	0.41
26:BD:3:GLY:HA3	26:BD:49:GLN:HE22	1.86	0.41
27:BE:6:LYS:HB2	27:BE:12:LEU:HD11	2.01	0.41
28:BF:32:LYS:HD2	28:BF:89:THR:HB	2.01	0.41
29:BG:97:VAL:HG21	29:BG:123:GLU:HA	2.02	0.41
52:BI:130:GLY:HA2	52:BI:133:ARG:NH2	2.35	0.41
31:BJ:102:GLU:HG2	31:BJ:103:ILE:N	2.35	0.41
34:BM:69:PRO:HB2	34:BM:70:ASP:H	1.62	0.41
35:BN:29:VAL:HG22	35:BN:83:LEU:CD2	2.48	0.41
35:BN:1:MET:SD	35:BN:2:ARG:HB2	2.60	0.41
37:BP:59:THR:O	37:BP:59:THR:HG22	2.20	0.41
38:BQ:80:ASN:HD22	38:BQ:80:ASN:HA	1.66	0.41
40:BS:85:ILE:HG22	40:BS:86:MET:N	2.35	0.41
24:BV:80:HIS:CD2	24:BV:81:PRO:HD2	2.56	0.41
23:BB:2270:A:C4'	43:BW:18:LYS:HB3	2.49	0.41
45:BY:8:GLN:NE2	45:BY:23:LEU:HD21	2.35	0.41
1:CA:1190:G:H5'	2:CC:175:HIS:NE2	2.36	0.41
1:CA:1293:C:H2'	1:CA:1294:G:O4'	2.20	0.41
1:CA:162:A:H2'	1:CA:163:C:O4'	2.20	0.41
1:CA:207:C:O2'	1:CA:208:U:H5'	2.20	0.41
1:CA:596:A:O2'	1:CA:597:G:H5'	2.20	0.41
1:CA:688:G:H5'	10:CK:48:GLY:CA	2.50	0.41
1:CA:821:G:H2'	1:CA:822:U:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:857:C:O2'	1:CA:858:G:H5'	2.21	0.41
1:CA:928:G:O2'	1:CA:929:G:H5'	2.20	0.41
20:CB:116:LEU:HD22	20:CB:140:LEU:CD2	2.41	0.41
20:CB:156:LEU:HD11	20:CB:178:LEU:CD1	2.33	0.41
2:CC:111:ASP:O	2:CC:112:ALA:HB3	2.21	0.41
4:CE:64:GLU:HA	4:CE:67:ARG:HH12	1.85	0.41
4:CE:89:THR:HG22	4:CE:91:SER:N	2.22	0.41
6:CG:24:LYS:HA	6:CG:27:ASN:HD22	1.85	0.41
9:CJ:44:THR:HG23	9:CJ:70:HIS:HA	2.02	0.41
11:CL:49:ARG:HD3	11:CL:89:LEU:HD21	2.01	0.41
15:CP:28:ARG:C	15:CP:30:GLY:H	2.24	0.41
17:CR:20:ILE:HD12	17:CR:53:GLN:NE2	2.34	0.41
10:CK:110:THR:HG22	21:CU:4:LYS:HA	2.02	0.41
48:D1:49:LYS:C	48:D1:49:LYS:HD2	2.41	0.41
50:D3:4:LYS:HZ1	50:D3:60:CYS:HB3	1.86	0.41
23:DB:1064:C:O2'	23:DB:1065:U:H5'	2.20	0.41
23:DB:1449:G:H2'	23:DB:1450:G:C8	2.56	0.41
23:DB:1460:U:H5''	23:DB:1461:C:C6	2.56	0.41
23:DB:1678:A:C2'	23:DB:1679:A:H5'	2.50	0.41
23:DB:1893:C:H2'	23:DB:1894:C:O4'	2.21	0.41
23:DB:2191:A:H2'	23:DB:2192:U:H6	1.83	0.41
23:DB:2221:G:H2'	23:DB:2222:C:C6	2.56	0.41
23:DB:2668:G:O2'	23:DB:2669:G:H5'	2.20	0.41
23:DB:2681:C:OP2	26:DD:114:LYS:HE3	2.20	0.41
23:DB:2686:G:H2'	23:DB:2687:U:C6	2.55	0.41
23:DB:413:C:O5'	23:DB:413:C:H6	2.04	0.41
23:DB:432:A:H2'	23:DB:433:C:O4'	2.20	0.41
23:DB:523:C:H4'	23:DB:540:C:O2	2.21	0.41
23:DB:570:G:C2'	23:DB:571:U:H5'	2.51	0.41
23:DB:63:A:OP2	23:DB:63:A:H8	2.04	0.41
23:DB:794:A:C4	23:DB:795:C:C5	3.09	0.41
23:DB:811:U:H2'	33:DL:32:GLY:HA2	2.02	0.41
23:DB:82:U:H2'	23:DB:83:A:O4'	2.20	0.41
23:DB:834:G:H2'	23:DB:835:C:C6	2.55	0.41
23:DB:921:C:C2	23:DB:922:C:C5	3.09	0.41
25:DC:107:LYS:HD2	25:DC:196:ASN:ND2	2.36	0.41
25:DC:59:GLN:HB2	25:DC:60:ALA:H	1.64	0.41
25:DC:28:PRO:CG	25:DC:79:ARG:HE	2.34	0.41
27:DE:15:SER:HB3	27:DE:196:VAL:HG13	2.02	0.41
28:DF:48:LEU:C	28:DF:50:ASP:H	2.24	0.41
29:DG:104:LEU:N	29:DG:112:VAL:O	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:DG:154:GLU:OE2	29:DG:159:LYS:HB2	2.21	0.41
30:DH:9:VAL:O	30:DH:11:ASN:N	2.54	0.41
52:DI:9:LYS:HG2	52:DI:57:VAL:HG22	2.03	0.41
31:DJ:84:ILE:CG1	31:DJ:85:LYS:H	2.33	0.41
32:DK:99:ILE:O	32:DK:119:ALA:HB2	2.20	0.41
34:DM:105:MET:O	34:DM:106:ASP:C	2.59	0.41
34:DM:119:LEU:CD2	34:DM:119:LEU:H	2.34	0.41
35:DN:38:LEU:O	35:DN:41:ALA:HB3	2.21	0.41
37:DP:28:LYS:HZ2	37:DP:44:GLY:H	1.67	0.41
24:DV:48:MET:O	24:DV:51:GLN:HG3	2.20	0.41
45:DY:26:LEU:HD23	45:DY:26:LEU:N	2.35	0.41
46:DZ:17:SER:H	46:DZ:21:VAL:HG12	1.85	0.41
46:DZ:54:GLY:N	46:DZ:57:VAL:HG23	2.15	0.41
1:AA:1010:U:H2'	1:AA:1011:C:H6	1.86	0.41
1:AA:1207:G:O2'	1:AA:1208:C:H5'	2.20	0.41
1:AA:1221:G:O2'	1:AA:1222:G:H5'	2.20	0.41
1:AA:1505:G:H5''	1:AA:1506:U:O5'	2.21	0.41
1:AA:102:G:N3	1:AA:151:A:H2	2.19	0.41
1:AA:191:G:H2'	1:AA:192:A:C8	2.56	0.41
1:AA:191:G:H2'	1:AA:192:A:H8	1.86	0.41
1:AA:42:G:O2'	1:AA:43:C:H5'	2.20	0.41
1:AA:619:U:C2	3:AD:131:ILE:HD12	2.56	0.41
1:AA:857:C:O2'	1:AA:858:G:H5'	2.21	0.41
1:AA:967:C:OP1	1:AA:969:A:C8	2.74	0.41
2:AC:17:TRP:C	2:AC:19:SER:H	2.23	0.41
2:AC:38:VAL:CG2	2:AC:39:ARG:N	2.83	0.41
4:AE:89:THR:HG21	4:AE:134:ASN:ND2	2.36	0.41
4:AE:35:LEU:HD21	4:AE:47:PHE:HB2	2.02	0.41
8:AI:12:LYS:N	8:AI:109:GLN:HE22	2.13	0.41
10:AK:88:PRO:HA	10:AK:92:ARG:CD	2.51	0.41
11:AL:113:ARG:HA	11:AL:118:VAL:CG2	2.51	0.41
11:AL:5:GLN:HA	11:AL:8:ARG:NE	2.35	0.41
12:AM:12:LYS:CB	12:AM:17:ALA:HB2	2.51	0.41
12:AM:12:LYS:HB3	12:AM:17:ALA:HB2	2.02	0.41
12:AM:21:ILE:C	12:AM:23:GLY:H	2.24	0.41
14:AO:15:GLY:C	14:AO:17:ASP:H	2.24	0.41
14:AO:63:ARG:NH1	14:AO:63:ARG:HB3	2.35	0.41
16:AQ:56:ASP:C	16:AQ:79:GLU:HB3	2.40	0.41
23:BB:2285:C:H41	48:B1:24:LYS:HZ2	1.69	0.41
23:BB:834:G:O2'	50:B3:50:SER:HB3	2.21	0.41
51:B4:2:LYS:HZ2	51:B4:2:LYS:HB3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:989:G:OP1	23:BB:1157:G:H4'	2.20	0.41
23:BB:1196:C:H2'	23:BB:1197:G:C8	2.56	0.41
23:BB:1434:A:OP1	23:BB:1434:A:H4'	2.20	0.41
23:BB:1827:U:C2'	23:BB:1828:G:H5'	2.51	0.41
23:BB:2094:A:H5'	30:BH:25:TYR:CD1	2.54	0.41
23:BB:2096:C:H2'	23:BB:2097:A:H8	1.86	0.41
23:BB:2467:C:H1'	34:BM:118:LYS:HG2	2.01	0.41
23:BB:2653:U:H2'	23:BB:2654:A:C8	2.55	0.41
23:BB:689:A:H2'	23:BB:690:G:H8	1.85	0.41
23:BB:844:A:C2	23:BB:845:A:N1	2.89	0.41
23:BB:967:U:C2	23:BB:968:C:C5	3.09	0.41
25:BC:81:GLU:O	25:BC:90:ILE:HG22	2.21	0.41
25:BC:90:ILE:HG23	25:BC:91:ALA:N	2.35	0.41
23:BB:2512:C:H5''	26:BD:125:TRP:CZ3	2.56	0.41
23:BB:1656:C:OP1	26:BD:141:ARG:HD3	2.21	0.41
26:BD:165:MET:CG	26:BD:166:GLY:H	2.33	0.41
27:BE:132:LYS:H	27:BE:132:LYS:HG3	1.55	0.41
27:BE:45:ALA:HB3	27:BE:89:PRO:HA	2.02	0.41
28:BF:36:ASN:HB2	28:BF:87:LYS:CA	2.33	0.41
30:BH:141:LYS:HE2	30:BH:141:LYS:HB2	1.86	0.41
23:BB:558:U:C5'	31:BJ:114:LEU:HD22	2.51	0.41
31:BJ:34:ARG:CG	31:BJ:39:LYS:HD2	2.50	0.41
33:BL:4:ASN:O	33:BL:6:LEU:HD22	2.21	0.41
34:BM:105:MET:HE3	34:BM:106:ASP:C	2.41	0.41
36:BO:103:VAL:HG23	36:BO:104:GLN:N	2.35	0.41
36:BO:30:ARG:O	36:BO:31:THR:O	2.39	0.41
36:BO:60:GLU:HG3	36:BO:61:GLN:N	2.35	0.41
36:BO:64:TYR:HD1	36:BO:74:VAL:HG21	1.84	0.41
26:BD:11:MET:CE	37:BP:9:GLN:HG2	2.51	0.41
40:BS:6:LYS:CB	40:BS:104:THR:HG22	2.50	0.41
42:BU:60:LYS:CG	42:BU:61:GLU:N	2.84	0.41
42:BU:6:ARG:C	42:BU:8:ASP:N	2.73	0.41
42:BU:82:VAL:HG12	42:BU:83:GLY:N	2.35	0.41
43:BW:54:ARG:HH11	43:BW:54:ARG:CG	2.31	0.41
23:BB:968:C:H4'	45:BY:16:LEU:HD23	2.03	0.41
46:BZ:21:VAL:HG12	46:BZ:23:LYS:HG2	2.03	0.41
1:CA:1043:G:C2'	1:CA:1044:A:O5'	2.68	0.41
1:CA:1128:C:H2'	1:CA:1129:C:O4'	2.21	0.41
1:CA:1170:A:H5''	20:CB:138:ARG:HH22	1.84	0.41
1:CA:1294:G:H2'	1:CA:1295:U:C6	2.56	0.41
1:CA:1510:C:H2'	1:CA:1511:G:C8	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:191:G:H2'	1:CA:192:A:C8	2.56	0.41
1:CA:844:G:H2'	1:CA:845:A:H5''	2.01	0.41
20:CB:173:LYS:HD2	20:CB:173:LYS:C	2.41	0.41
3:CD:115:GLN:HE21	3:CD:153:ARG:HH12	1.66	0.41
6:CG:70:PRO:HD3	6:CG:102:TRP:CZ3	2.56	0.41
8:CI:10:ARG:HA	8:CI:14:SER:O	2.21	0.41
9:CJ:92:LEU:O	9:CJ:93:ALA:HB2	2.21	0.41
12:CM:13:HIS:HB2	12:CM:16:ILE:HG22	2.02	0.41
13:CN:12:ARG:HD3	13:CN:58:ARG:O	2.20	0.41
16:CQ:64:ARG:HH11	16:CQ:64:ARG:CB	2.34	0.41
18:CS:14:LEU:HD12	18:CS:14:LEU:C	2.41	0.41
18:CS:28:LYS:HB2	18:CS:29:PRO:HD2	2.02	0.41
19:CT:11:ILE:HG13	19:CT:12:GLN:N	2.36	0.41
19:CT:24:ARG:HG3	19:CT:65:LEU:HD11	2.03	0.41
48:D1:13:SER:HB3	48:D1:49:LYS:HZ1	1.84	0.41
48:D1:8:ILE:HG23	48:D1:9:LYS:N	2.35	0.41
50:D3:38:LYS:N	50:D3:38:LYS:HD2	2.36	0.41
23:DB:1322:A:C5	23:DB:1323:C:C5	3.08	0.41
23:DB:1446:C:H2'	23:DB:1447:C:H6	1.86	0.41
23:DB:1498:C:H2'	23:DB:1499:C:H6	1.86	0.41
23:DB:1744:A:O2'	23:DB:1745:A:H5'	2.21	0.41
23:DB:2070:A:C2	23:DB:2442:C:C2	3.09	0.41
23:DB:2259:U:C2	23:DB:2260:C:C6	3.09	0.41
23:DB:31:C:C2'	23:DB:32:C:H5'	2.51	0.41
23:DB:484:C:H2'	23:DB:485:C:C6	2.54	0.41
23:DB:622:G:H2'	23:DB:623:C:H6	1.86	0.41
23:DB:67:U:H2'	23:DB:68:G:O4'	2.21	0.41
25:DC:265:PHE:HB3	25:DC:266:ILE:H	1.75	0.41
26:DD:160:LYS:O	26:DD:161:MET:C	2.58	0.41
28:DF:35:LEU:CD2	28:DF:153:ILE:HG12	2.51	0.41
29:DG:121:THR:HG22	29:DG:122:ALA:H	1.86	0.41
29:DG:145:ALA:HB2	29:DG:163:TYR:OH	2.20	0.41
29:DG:156:TYR:HD1	29:DG:171:LYS:HG2	1.85	0.41
29:DG:2:ARG:HD2	29:DG:2:ARG:O	2.21	0.41
30:DH:3:VAL:CB	30:DH:37:VAL:HG11	2.50	0.41
52:DI:131:THR:O	52:DI:135:MET:HG3	2.21	0.41
33:DL:66:PHE:CD2	33:DL:67:THR:N	2.89	0.41
35:DN:14:SER:HA	35:DN:17:ARG:CZ	2.51	0.41
37:DP:91:VAL:O	37:DP:92:ARG:HB3	2.21	0.41
39:DR:20:VAL:HA	39:DR:97:LYS:O	2.20	0.41
39:DR:35:PHE:O	39:DR:63:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:61:LEU:O	41:DT:62:VAL:HG13	2.21	0.41
1:AA:1005:A:C2'	1:AA:1006:G:H5'	2.51	0.41
1:AA:1148:U:C2'	1:AA:1149:C:H5'	2.50	0.41
1:AA:1245:C:O2'	1:AA:1246:A:H5'	2.20	0.41
1:AA:144:G:H2'	1:AA:145:G:C8	2.56	0.41
1:AA:1523:G:H2'	1:AA:1524:C:C6	2.55	0.41
1:AA:208:U:H2'	1:AA:210:C:C6	2.55	0.41
1:AA:401:C:H2'	1:AA:402:G:H8	1.86	0.41
1:AA:493:A:N3	1:AA:493:A:O4'	2.54	0.41
1:AA:688:G:O2'	1:AA:689:C:H5'	2.20	0.41
1:AA:737:C:H2'	1:AA:738:C:C6	2.55	0.41
1:AA:955:U:H2'	1:AA:956:U:O4'	2.21	0.41
20:AB:114:LYS:HA	20:AB:117:GLU:OE1	2.21	0.41
2:AC:139:ASN:ND2	2:AC:139:ASN:N	2.68	0.41
2:AC:178:ARG:O	2:AC:206:ILE:HA	2.21	0.41
3:AD:89:LEU:O	3:AD:92:LEU:HB2	2.21	0.41
5:AF:38:ARG:HH21	5:AF:96:VAL:CG1	2.33	0.41
6:AG:94:ARG:NH1	6:AG:98:LEU:HD11	2.35	0.41
7:AH:9:MET:O	7:AH:13:ILE:HG13	2.20	0.41
10:AK:122:PRO:HG2	21:AU:35:GLU:HG2	2.02	0.41
10:AK:57:SER:O	10:AK:58:THR:C	2.59	0.41
12:AM:94:LEU:HD22	12:AM:101:THR:OG1	2.21	0.41
14:AO:71:ARG:HB2	14:AO:71:ARG:NH1	2.35	0.41
16:AQ:8:GLN:HA	16:AQ:58:VAL:O	2.21	0.41
19:AT:20:ASN:HA	19:AT:65:LEU:HD21	2.02	0.41
48:B1:29:LYS:CE	48:B1:30:PRO:HD2	2.47	0.41
22:BA:86:G:H3'	22:BA:87:U:H6	1.84	0.41
22:BA:95:U:H2'	22:BA:96:G:C8	2.55	0.41
23:BB:1460:U:H5''	23:BB:1461:C:C6	2.56	0.41
23:BB:2019:A:O2'	38:BQ:32:ARG:NH2	2.54	0.41
23:BB:2026:U:H2'	23:BB:2027:G:C8	2.55	0.41
23:BB:2545:G:N3	23:BB:2565:A:H2	2.19	0.41
23:BB:2897:U:H2'	23:BB:2898:U:C6	2.55	0.41
23:BB:356:G:O2'	23:BB:357:C:H5'	2.20	0.41
23:BB:503:A:H5''	23:BB:504:A:H5'	2.02	0.41
23:BB:607:U:C4	23:BB:620:G:O4'	2.73	0.41
23:BB:723:C:H2'	23:BB:724:U:H6	1.84	0.41
23:BB:794:A:C4	23:BB:795:C:C5	3.09	0.41
23:BB:849:A:H2'	23:BB:850:U:C6	2.55	0.41
25:BC:160:TYR:CD2	25:BC:193:GLU:HG2	2.56	0.41
25:BC:24:HIS:HB3	25:BC:27:LYS:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:BC:37:SER:H	25:BC:62:ARG:HB2	1.86	0.41
26:BD:140:HIS:O	26:BD:142:VAL:N	2.54	0.41
26:BD:145:SER:C	26:BD:146:ILE:HD12	2.40	0.41
26:BD:170:VAL:CG1	26:BD:194:PRO:HG2	2.51	0.41
27:BE:190:ALA:HB1	27:BE:193:VAL:CB	2.50	0.41
27:BE:25:GLU:O	27:BE:27:LEU:N	2.54	0.41
30:BH:69:ALA:HB1	30:BH:141:LYS:HZ3	1.85	0.41
30:BH:41:LYS:N	30:BH:41:LYS:HD2	2.33	0.41
52:BI:16:MET:N	52:BI:42:ASN:OD1	2.54	0.41
31:BJ:125:TYR:CE2	31:BJ:134:ALA:HB2	2.56	0.41
31:BJ:84:ILE:CG2	31:BJ:85:LYS:H	2.31	0.41
32:BK:91:SER:O	32:BK:92:GLU:C	2.59	0.41
33:BL:118:THR:OG1	33:BL:119:PRO:HD3	2.20	0.41
33:BL:84:LYS:O	33:BL:85:VAL:CB	2.69	0.41
34:BM:45:GLN:CG	34:BM:120:ALA:HA	2.51	0.41
36:BO:105:ALA:HA	36:BO:108:ASP:OD2	2.21	0.41
37:BP:45:VAL:HG12	37:BP:65:ASN:HB2	2.02	0.41
38:BQ:108:LEU:HD21	39:BR:52:PRO:O	2.21	0.41
42:BU:9:GLU:CG	42:BU:23:LYS:HB3	2.51	0.41
1:CA:1039:G:O2'	1:CA:1040:U:H5'	2.21	0.41
1:CA:359:G:O2'	1:CA:360:G:H5'	2.21	0.41
1:CA:746:A:H2'	1:CA:747:A:H8	1.86	0.41
1:CA:90:C:H2'	1:CA:91:U:H6	1.84	0.41
1:CA:984:C:H2'	1:CA:985:C:C6	2.56	0.41
4:CE:109:ALA:HB1	4:CE:136:VAL:CG1	2.49	0.41
4:CE:14:LEU:HD13	4:CE:14:LEU:C	2.41	0.41
7:CH:100:ILE:HD12	7:CH:101:ALA:N	2.36	0.41
12:CM:48:SER:C	12:CM:50:GLY:H	2.23	0.41
16:CQ:57:VAL:HB	16:CQ:79:GLU:HB2	2.03	0.41
1:CA:1317:C:H42	18:CS:36:ARG:NH2	2.18	0.41
21:CU:35:GLU:OE1	21:CU:35:GLU:HA	2.20	0.41
50:D3:32:LEU:HA	50:D3:32:LEU:HD22	1.93	0.41
51:D4:36:ARG:HG2	51:D4:36:ARG:NH1	2.34	0.41
22:DA:92:C:H2'	22:DA:93:C:C6	2.56	0.41
23:DB:1268:A:H2'	23:DB:1269:A:C8	2.55	0.41
23:DB:1655:A:H2'	23:DB:1656:C:O4'	2.21	0.41
23:DB:1825:U:H5'	25:DC:244:VAL:HG23	1.99	0.41
23:DB:1897:G:O2'	23:DB:1898:U:H5'	2.20	0.41
23:DB:2597:G:OP1	25:DC:239:PHE:CD2	2.74	0.41
23:DB:2665:A:H2'	23:DB:2666:C:O2	2.21	0.41
23:DB:2710:C:H2'	23:DB:2711:A:C8	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:619:G:H3'	23:DB:620:G:H21	1.86	0.41
23:DB:820:A:H2'	23:DB:821:A:O4'	2.20	0.41
23:DB:869:G:H2'	23:DB:870:U:O4'	2.20	0.41
25:DC:163:ILE:HA	25:DC:173:LEU:HD23	2.03	0.41
25:DC:173:LEU:HD12	25:DC:183:VAL:CG1	2.48	0.41
26:DD:26:VAL:HG13	26:DD:26:VAL:O	2.21	0.41
27:DE:48:THR:CG2	27:DE:85:PHE:N	2.84	0.41
28:DF:99:PHE:CD1	28:DF:102:LEU:HD23	2.56	0.41
31:DJ:128:ASN:O	31:DJ:130:HIS:N	2.54	0.41
31:DJ:36:LEU:O	31:DJ:51:GLY:HA2	2.21	0.41
33:DL:33:ARG:HB3	39:DR:85:LYS:NZ	2.35	0.41
35:DN:106:ASP:O	35:DN:107:ASN:HB3	2.20	0.41
36:DO:10:ARG:HD2	36:DO:94:ARG:HG3	2.02	0.41
40:DS:20:VAL:O	40:DS:24:ILE:HG13	2.21	0.41
42:DU:71:ILE:CG1	42:DU:72:PHE:N	2.84	0.41
24:DV:43:ASP:O	24:DV:47:VAL:HG23	2.20	0.41
43:DW:18:LYS:HG3	43:DW:19:ARG:N	2.36	0.41
44:DX:40:SER:OG	44:DX:42:LEU:HD23	2.20	0.41
44:DX:41:HIS:ND1	44:DX:43:LEU:HD23	2.36	0.41
46:DZ:1:MET:HG3	46:DZ:9:TYR:CE2	2.56	0.41
46:DZ:20:ASN:O	46:DZ:21:VAL:CB	2.63	0.41
46:DZ:11:GLU:O	46:DZ:27:THR:HG22	2.21	0.41
1:AA:1033:G:C3'	1:AA:1034:G:H5''	2.51	0.41
1:AA:1258:G:O2'	1:AA:1259:C:H5'	2.21	0.41
1:AA:1458:G:H2'	1:AA:1459:G:C8	2.56	0.41
1:AA:31:G:H2'	1:AA:48:C:H41	1.86	0.41
1:AA:601:G:O2'	1:AA:602:A:H5'	2.21	0.41
1:AA:846:G:H5'	1:AA:847:G:OP2	2.21	0.41
1:AA:890:G:O2'	1:AA:906:A:N6	2.53	0.41
1:AA:91:U:H2'	1:AA:92:U:C6	2.56	0.41
1:AA:960:U:C1'	1:AA:961:U:OP2	2.69	0.41
20:AB:202:ASN:ND2	20:AB:204:ASP:N	2.60	0.41
20:AB:96:LEU:HD23	20:AB:99:MET:CE	2.50	0.41
2:AC:137:VAL:HG11	2:AC:150:VAL:HG21	2.01	0.41
2:AC:176:THR:O	2:AC:179:ALA:HB3	2.21	0.41
1:AA:1057:G:H4'	2:AC:194:VAL:O	2.21	0.41
3:AD:37:PRO:HD2	3:AD:41:GLY:CA	2.51	0.41
3:AD:55:ARG:NH2	3:AD:58:GLN:NE2	2.67	0.41
5:AF:7:VAL:HB	5:AF:61:LEU:HD22	2.03	0.41
1:AA:736:C:H4'	5:AF:88:MET:HE3	2.02	0.41
7:AH:113:ARG:HE	7:AH:113:ARG:C	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:35:GLU:N	8:AI:35:GLU:OE2	2.54	0.41
8:AI:56:MET:HA	8:AI:59:LYS:HB2	2.03	0.41
9:AJ:67:ILE:CG1	13:AN:95:LEU:HD13	2.44	0.41
14:AO:15:GLY:HA3	14:AO:20:ASP:OD1	2.21	0.41
14:AO:52:ARG:HG3	14:AO:55:LEU:HD23	2.03	0.41
17:AR:24:ASP:O	17:AR:28:LEU:HD13	2.20	0.41
49:B2:9:VAL:O	49:B2:12:ARG:HB2	2.20	0.41
51:B4:11:CYS:SG	51:B4:12:ARG:N	2.94	0.41
51:B4:24:ARG:NH2	51:B4:36:ARG:HH11	2.19	0.41
23:BB:1142:A:C4	23:BB:1144:A:C8	3.08	0.41
23:BB:1192:G:C2'	23:BB:1193:G:H5'	2.50	0.41
23:BB:134:G:C6	23:BB:146:A:C6	3.08	0.41
23:BB:137:U:H2'	23:BB:138:U:C5'	2.51	0.41
23:BB:1568:G:H5''	25:BC:60:ALA:HB3	2.02	0.41
23:BB:1764:C:H2'	23:BB:1765:U:C6	2.56	0.41
23:BB:1795:C:H2'	23:BB:1796:U:H6	1.85	0.41
23:BB:1989:G:H2'	23:BB:1990:C:O4'	2.21	0.41
23:BB:2441:U:O2'	23:BB:2442:C:H5'	2.21	0.41
23:BB:2617:U:O2'	23:BB:2618:G:H5'	2.21	0.41
23:BB:2653:U:HO2'	29:BG:109:SER:CB	2.34	0.41
23:BB:2720:U:H2'	23:BB:2721:A:C8	2.55	0.41
23:BB:286:U:H2'	23:BB:287:G:H8	1.86	0.41
25:BC:107:LYS:H	25:BC:194:VAL:CG2	2.31	0.41
25:BC:185:ALA:C	25:BC:187:CYS:H	2.24	0.41
25:BC:143:VAL:HG12	25:BC:187:CYS:SG	2.61	0.41
26:BD:46:ARG:HD2	26:BD:80:TRP:CH2	2.55	0.41
27:BE:146:VAL:HG23	27:BE:148:ILE:HG23	2.02	0.41
27:BE:153:LEU:HD11	27:BE:193:VAL:HG11	2.02	0.41
30:BH:66:ASN:OD1	30:BH:138:VAL:HG11	2.20	0.41
31:BJ:35:ARG:C	31:BJ:37:ARG:H	2.24	0.41
32:BK:108:ARG:HG3	32:BK:108:ARG:NH1	2.36	0.41
23:BB:196:A:N7	33:BL:52:GLY:HA2	2.36	0.41
34:BM:8:LYS:NZ	34:BM:69:PRO:HG2	2.36	0.41
23:BB:2020:A:OP1	38:BQ:25:GLY:HA3	2.20	0.41
23:BB:1225:G:C5'	39:BR:78:ARG:NH2	2.84	0.41
42:BU:71:ILE:HG21	42:BU:95:PHE:CE1	2.55	0.41
24:BV:87:GLN:N	24:BV:87:GLN:NE2	2.69	0.41
43:BW:81:ILE:HG13	43:BW:82:GLU:HG2	2.03	0.41
45:BY:16:LEU:CB	45:BY:17:PRO:CD	2.91	0.41
46:BZ:3:LYS:HG3	46:BZ:6:HIS:O	2.21	0.41
1:CA:181:A:H1'	1:CA:182:A:C2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:599:C:H2'	1:CA:600:A:H8	1.85	0.41
1:CA:607:A:H2'	1:CA:608:A:H8	1.86	0.41
1:CA:672:U:H2'	1:CA:673:A:H8	1.83	0.41
1:CA:728:A:H2'	1:CA:729:A:C8	2.55	0.41
1:CA:737:C:H2'	1:CA:738:C:C6	2.55	0.41
1:CA:738:C:H2'	1:CA:739:C:H6	1.86	0.41
1:CA:828:U:H2'	1:CA:829:G:O5'	2.21	0.41
1:CA:830:G:H2'	1:CA:831:A:H8	1.86	0.41
1:CA:407:U:O2'	3:CD:112:GLU:HG3	2.21	0.41
3:CD:160:LEU:C	3:CD:160:LEU:HD13	2.41	0.41
8:CI:109:GLN:HE21	8:CI:109:GLN:HB3	1.61	0.41
8:CI:56:MET:C	8:CI:58:GLU:N	2.73	0.41
8:CI:20:ILE:CG2	8:CI:60:LEU:HD22	2.51	0.41
11:CL:29:LYS:HB3	11:CL:56:LEU:HD22	2.02	0.41
11:CL:79:ILE:C	11:CL:101:LEU:HD12	2.41	0.41
11:CL:49:ARG:CD	11:CL:89:LEU:HD21	2.51	0.41
1:CA:750:C:O2	14:CO:22:GLY:HA3	2.21	0.41
14:CO:49:HIS:O	14:CO:52:ARG:HB3	2.21	0.41
18:CS:41:PRO:HA	18:CS:44:ILE:CG1	2.50	0.41
48:D1:34:GLU:HB3	48:D1:50:GLU:HB3	2.03	0.41
48:D1:5:ARG:HG2	48:D1:5:ARG:NH1	2.35	0.41
23:DB:247:G:O6	50:D3:7:ARG:HD2	2.21	0.41
51:D4:15:LYS:C	51:D4:17:VAL:N	2.74	0.41
51:D4:23:ILE:CD1	51:D4:24:ARG:HH11	2.34	0.41
23:DB:103:A:H3'	23:DB:104:A:H8	1.85	0.41
23:DB:1099:G:C5'	52:DI:3:LYS:CA	2.99	0.41
23:DB:1120:G:H2'	23:DB:1121:C:C6	2.56	0.41
23:DB:1390:U:O2'	23:DB:1391:U:H5'	2.20	0.41
23:DB:1528:A:C2'	23:DB:1529:G:H5'	2.50	0.41
23:DB:1593:A:H2'	23:DB:1594:U:O4'	2.20	0.41
23:DB:1841:U:H2'	23:DB:1842:G:C8	2.55	0.41
23:DB:1924:C:O2'	23:DB:1925:C:H5'	2.21	0.41
23:DB:2188:U:O2'	23:DB:2189:U:H5'	2.21	0.41
23:DB:2559:C:O2'	23:DB:2560:A:H5'	2.19	0.41
23:DB:296:U:H3'	55:DB:3414:HOH:O	2.21	0.41
23:DB:301:G:C4	23:DB:302:C:C6	3.09	0.41
23:DB:583:G:H2'	23:DB:584:C:H6	1.86	0.41
25:DC:103:ILE:HG22	25:DC:104:LEU:N	2.36	0.41
26:DD:100:LEU:HD13	26:DD:100:LEU:C	2.41	0.41
26:DD:197:THR:CG2	26:DD:198:GLY:H	2.08	0.41
23:DB:443:A:C8	27:DE:40:ARG:HG2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:DF:98:PHE:HA	28:DF:101:ARG:CG	2.50	0.41
29:DG:152:ARG:HA	29:DG:153:PRO:HD3	1.96	0.41
30:DH:67:ALA:C	30:DH:69:ALA:H	2.23	0.41
52:DI:100:ILE:CG2	52:DI:104:GLN:HB2	2.50	0.41
34:DM:127:LYS:C	34:DM:127:LYS:HD3	2.41	0.41
36:DO:30:ARG:CD	36:DO:97:PHE:HD2	2.33	0.41
37:DP:61:ARG:CB	37:DP:61:ARG:CZ	2.98	0.41
37:DP:91:VAL:O	37:DP:92:ARG:CB	2.69	0.41
42:DU:26:ASN:O	42:DU:28:LEU:HG	2.21	0.41
44:DX:7:ARG:C	44:DX:8:GLU:HG3	2.41	0.41
45:DY:24:LEU:C	45:DY:26:LEU:N	2.74	0.41
45:DY:5:LYS:NZ	45:DY:5:LYS:CB	2.84	0.41
1:AA:1276:G:H2'	1:AA:1277:C:C6	2.56	0.41
1:AA:1401:G:O2'	1:AA:1402:C:H5'	2.20	0.41
1:AA:358:U:H2'	1:AA:359:G:H8	1.86	0.41
1:AA:359:G:O2'	1:AA:360:G:H5'	2.21	0.41
1:AA:393:A:O2'	1:AA:394:G:H5'	2.20	0.41
1:AA:367:U:OP1	1:AA:395:C:H1'	2.21	0.41
1:AA:580:C:H2'	1:AA:581:G:O4'	2.21	0.41
1:AA:611:C:H2'	1:AA:612:C:C6	2.56	0.41
1:AA:612:C:H2'	1:AA:613:C:C6	2.56	0.41
20:AB:140:LEU:HD22	20:AB:144:GLU:OE2	2.20	0.41
3:AD:197:HIS:HA	3:AD:200:VAL:HG22	2.03	0.41
4:AE:81:GLN:NE2	4:AE:82:HIS:ND1	2.68	0.41
8:AI:9:GLY:O	8:AI:16:ALA:HB3	2.21	0.41
8:AI:52:GLU:N	8:AI:56:MET:HG2	2.35	0.41
13:AN:68:ARG:HH11	13:AN:70:HIS:HB2	1.85	0.41
16:AQ:10:ARG:HD3	16:AQ:55:GLY:H	1.86	0.41
48:B1:11:VAL:HG13	48:B1:12:SER:N	2.35	0.41
49:B2:38:GLY:O	49:B2:39:ARG:C	2.58	0.41
49:B2:39:ARG:HB2	49:B2:39:ARG:NH1	2.35	0.41
50:B3:4:LYS:HZ1	50:B3:59:ALA:H	1.69	0.41
22:BA:109:A:O2'	22:BA:110:C:H5'	2.20	0.41
23:BB:1060:U:C1'	23:BB:1062:G:H5'	2.51	0.41
23:BB:1170:C:O2'	23:BB:1171:G:H5'	2.21	0.41
23:BB:1198:U:H2'	23:BB:1199:U:C5	2.55	0.41
23:BB:1449:G:H2'	23:BB:1450:G:C8	2.56	0.41
23:BB:1936:A:C2	23:BB:1943:U:H5	2.39	0.41
23:BB:197:A:H4'	23:BB:2069:G:OP2	2.21	0.41
23:BB:2139:U:O5'	23:BB:2139:U:H6	2.04	0.41
23:BB:2621:G:P	26:BD:124:ARG:HH22	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:264:C:C2'	23:BB:265:A:H5''	2.51	0.41
23:BB:2693:G:O2'	23:BB:2694:G:H5'	2.21	0.41
23:BB:527:C:H5''	23:BB:2779:U:N3	2.36	0.41
23:BB:28:A:H2'	23:BB:29:U:H6	1.85	0.41
23:BB:570:G:C2'	23:BB:571:U:H5'	2.51	0.41
23:BB:730:A:C2	23:BB:731:C:C6	3.09	0.41
23:BB:820:A:H2'	23:BB:821:A:O4'	2.21	0.41
25:BC:152:GLN:HB3	25:BC:153:LEU:H	1.53	0.41
25:BC:205:GLY:O	25:BC:206:LYS:HG2	2.21	0.41
28:BF:116:LEU:HG	28:BF:175:PRO:O	2.21	0.41
30:BH:41:LYS:O	30:BH:42:LYS:C	2.59	0.41
31:BJ:16:TYR:HD2	31:BJ:140:LEU:HB2	1.86	0.41
33:BL:2:ARG:NH1	33:BL:4:ASN:HD21	2.19	0.41
34:BM:8:LYS:HZ2	34:BM:69:PRO:HB2	1.84	0.41
34:BM:8:LYS:NZ	34:BM:69:PRO:HB2	2.36	0.41
35:BN:103:ARG:HG2	35:BN:103:ARG:O	2.21	0.41
36:BO:36:TYR:HA	36:BO:52:SER:HB2	2.03	0.41
37:BP:28:LYS:HB3	37:BP:44:GLY:H	1.85	0.41
37:BP:5:LYS:C	37:BP:7:LEU:N	2.71	0.41
38:BQ:50:ARG:HG2	38:BQ:50:ARG:HH11	1.86	0.41
39:BR:73:LYS:O	39:BR:74:ILE:O	2.39	0.41
40:BS:37:THR:HG23	40:BS:38:TYR:CD1	2.56	0.41
42:BU:38:ILE:HB	42:BU:62:ALA:HB3	2.03	0.41
42:BU:41:VAL:O	42:BU:41:VAL:HG13	2.20	0.41
43:BW:14:ASP:O	43:BW:15:SER:HB3	2.21	0.41
44:BX:50:VAL:O	44:BX:53:VAL:HG23	2.21	0.41
1:CA:1004:A:H3'	1:CA:1005:A:C8	2.54	0.41
1:CA:1015:G:O2'	1:CA:1016:A:H5'	2.21	0.41
1:CA:1018:G:H2'	1:CA:1019:A:C8	2.55	0.41
1:CA:1029:U:N3	1:CA:1033:G:C2	2.89	0.41
1:CA:175:C:H2'	1:CA:176:C:H6	1.86	0.41
1:CA:208:U:C2'	1:CA:209:U:H5''	2.50	0.41
1:CA:276:G:O2'	1:CA:277:C:H5'	2.21	0.41
1:CA:358:U:H2'	1:CA:359:G:H8	1.86	0.41
1:CA:493:A:O4'	1:CA:493:A:N3	2.54	0.41
1:CA:691:G:H1'	1:CA:696:A:H61	1.84	0.41
1:CA:878:A:H2'	1:CA:879:C:C6	2.56	0.41
20:CB:148:GLY:C	20:CB:150:ILE:H	2.24	0.41
2:CC:37:LYS:HB3	2:CC:41:TYR:CE1	2.56	0.41
2:CC:91:ALA:HB2	2:CC:98:ALA:CB	2.51	0.41
4:CE:21:SER:HA	4:CE:29:ILE:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CH:17:GLN:NE2	7:CH:69:ALA:HB1	2.36	0.41
8:CI:35:GLU:O	8:CI:39:GLY:HA2	2.21	0.41
8:CI:94:ARG:HH11	8:CI:98:ARG:NH1	2.19	0.41
13:CN:29:ILE:HB	13:CN:44:VAL:HG21	2.03	0.41
16:CQ:56:ASP:HA	16:CQ:81:ALA:HA	2.03	0.41
17:CR:47:ARG:HH12	17:CR:49:LYS:H	1.69	0.41
19:CT:38:ILE:HA	19:CT:85:LEU:HD22	2.02	0.41
22:DA:70:C:H2'	22:DA:71:C:H6	1.85	0.41
22:DA:76:G:O2'	22:DA:77:U:H5'	2.21	0.41
23:DB:1030:C:O2'	23:DB:1031:G:H5'	2.21	0.41
23:DB:1041:G:C2	23:DB:1042:G:N7	2.88	0.41
23:DB:1827:U:C2'	23:DB:1828:G:H5'	2.51	0.41
23:DB:1831:G:H2'	23:DB:1832:C:C6	2.56	0.41
23:DB:2195:U:O2'	23:DB:2196:C:H5'	2.21	0.41
23:DB:2308:G:N3	23:DB:2308:G:H3'	2.35	0.41
23:DB:2441:U:O2'	23:DB:2442:C:H5'	2.21	0.41
23:DB:266:G:O4'	23:DB:266:G:OP2	2.38	0.41
23:DB:2693:G:O2'	23:DB:2694:G:H5'	2.21	0.41
23:DB:2694:G:O2'	23:DB:2695:U:H5'	2.21	0.41
23:DB:2825:G:C2'	23:DB:2826:A:H5'	2.51	0.41
23:DB:288:U:H2'	23:DB:289:G:C8	2.56	0.41
23:DB:349:U:O2	23:DB:349:U:H2'	2.20	0.41
23:DB:458:G:O2'	23:DB:459:U:OP2	2.39	0.41
23:DB:927:A:H2'	23:DB:928:A:C8	2.55	0.41
25:DC:115:ILE:O	25:DC:116:GLN:HG3	2.21	0.41
25:DC:144:GLU:HG2	25:DC:150:GLY:HA2	2.02	0.41
25:DC:62:ARG:O	25:DC:62:ARG:NE	2.53	0.41
25:DC:63:ILE:HG23	25:DC:102:TYR:HD1	1.82	0.41
26:DD:160:LYS:HG3	26:DD:160:LYS:O	2.21	0.41
29:DG:145:ALA:HA	29:DG:148:ARG:HG2	2.02	0.41
29:DG:25:ILE:O	29:DG:31:GLU:HA	2.21	0.41
30:DH:68:ARG:HE	30:DH:71:LYS:CE	2.26	0.41
52:DI:8:VAL:HG11	52:DI:30:GLN:HG3	2.03	0.41
32:DK:63:VAL:HB	32:DK:83:ALA:CB	2.51	0.41
33:DL:25:SER:HB2	33:DL:26:GLY:H	1.55	0.41
34:DM:112:LEU:HA	34:DM:112:LEU:HD23	1.89	0.41
34:DM:124:LEU:O	34:DM:124:LEU:HD12	2.21	0.41
34:DM:16:ARG:NE	34:DM:18:ARG:NH1	2.65	0.41
34:DM:66:ARG:HA	34:DM:66:ARG:NE	2.35	0.41
35:DN:98:LEU:HD11	35:DN:114:GLU:OE2	2.20	0.41
37:DP:27:VAL:O	37:DP:29:VAL:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:27:VAL:HB	37:DP:86:LYS:O	2.22	0.41
38:DQ:102:LYS:O	38:DQ:106:THR:HG22	2.20	0.41
38:DQ:95:ALA:O	38:DQ:98:ALA:HB3	2.20	0.41
39:DR:10:LYS:O	39:DR:11:GLN:HB2	2.21	0.41
42:DU:95:PHE:HB2	42:DU:99:SER:C	2.41	0.41
44:DX:31:GLN:NE2	44:DX:31:GLN:CA	2.83	0.41
1:AA:1015:G:O2'	1:AA:1016:A:H5'	2.20	0.40
1:AA:1058:G:H2'	1:AA:1059:C:O4'	2.20	0.40
1:AA:1072:G:H2'	1:AA:1073:U:H6	1.86	0.40
1:AA:1150:A:N6	1:AA:1151:A:N6	2.69	0.40
1:AA:1219:A:OP1	13:AN:52:ARG:HG2	2.20	0.40
1:AA:1309:G:O2'	1:AA:1310:G:H5'	2.20	0.40
1:AA:1386:G:H2'	1:AA:1387:G:H8	1.85	0.40
1:AA:1462:C:H2'	1:AA:1463:U:H6	1.86	0.40
1:AA:213:G:C8	1:AA:214:C:C5	3.08	0.40
1:AA:276:G:O2'	1:AA:277:C:H5'	2.21	0.40
1:AA:37:U:O2'	1:AA:38:G:H5'	2.21	0.40
1:AA:415:A:O4'	1:AA:415:A:N3	2.53	0.40
1:AA:80:A:C2	1:AA:81:A:C4	3.09	0.40
1:AA:951:G:O2'	1:AA:952:U:H5'	2.21	0.40
20:AB:88:GLN:HE21	20:AB:88:GLN:HB2	1.65	0.40
2:AC:110:LEU:HD22	2:AC:145:ALA:HB2	2.04	0.40
2:AC:132:ALA:HA	2:AC:136:ALA:HB3	2.02	0.40
2:AC:40:GLN:CD	2:AC:44:LYS:HD2	2.42	0.40
4:AE:17:VAL:CG2	4:AE:34:ALA:HB2	2.50	0.40
6:AG:101:ARG:O	6:AG:105:GLU:N	2.46	0.40
7:AH:47:ASP:CG	7:AH:48:PHE:H	2.25	0.40
9:AJ:48:ARG:HA	9:AJ:66:GLU:HA	2.03	0.40
11:AL:56:LEU:H	11:AL:56:LEU:CD1	2.34	0.40
13:AN:27:LYS:NZ	13:AN:28:ALA:HB2	2.36	0.40
13:AN:92:ILE:HG21	13:AN:95:LEU:HD23	2.03	0.40
14:AO:34:GLN:HG2	14:AO:58:MET:HE1	2.03	0.40
16:AQ:64:ARG:HG3	16:AQ:65:PRO:HD2	2.03	0.40
18:AS:62:THR:HG22	18:AS:63:ASP:N	2.36	0.40
47:B0:39:ARG:HG2	47:B0:39:ARG:NH1	2.36	0.40
23:BB:2347:C:OP2	50:B3:37:THR:HG21	2.21	0.40
23:BB:2479:U:OP1	51:B4:1:MET:HE1	2.22	0.40
23:BB:330:A:N7	23:BB:1210:G:H2'	2.37	0.40
23:BB:1446:C:H2'	23:BB:1447:C:H6	1.86	0.40
23:BB:1537:G:C2'	23:BB:1538:G:H5'	2.51	0.40
23:BB:1765:U:O2'	23:BB:1766:G:H5'	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:175:G:O2'	23:BB:176:A:H5'	2.21	0.40
23:BB:1998:A:OP2	26:BD:141:ARG:NH2	2.54	0.40
23:BB:2221:G:H2'	23:BB:2222:C:C6	2.57	0.40
23:BB:2285:C:N4	48:B1:24:LYS:NZ	2.69	0.40
23:BB:2408:U:H2'	23:BB:2409:G:C8	2.55	0.40
23:BB:2595:G:N2	23:BB:2597:G:H3'	2.37	0.40
23:BB:2756:U:H5'	51:B4:18:LYS:NZ	2.36	0.40
23:BB:298:G:H1'	23:BB:340:A:H61	1.86	0.40
23:BB:600:G:H2'	23:BB:601:C:H6	1.84	0.40
23:BB:619:G:H3'	23:BB:620:G:H21	1.86	0.40
23:BB:949:G:O2'	23:BB:950:G:H5'	2.21	0.40
25:BC:229:HIS:O	25:BC:231:HIS:N	2.40	0.40
23:BB:691:C:H1'	25:BC:42:ARG:NH1	2.36	0.40
1:AA:807:A:O2'	25:BC:4:LYS:HB2	2.21	0.40
25:BC:67:LYS:O	25:BC:68:ARG:CB	2.66	0.40
26:BD:170:VAL:CG1	26:BD:171:THR:H	2.13	0.40
26:BD:68:PHE:O	26:BD:70:LYS:N	2.46	0.40
23:BB:2306:C:H1'	28:BF:132:ARG:NH2	2.37	0.40
29:BG:6:ALA:H	29:BG:7:PRO:CD	2.34	0.40
30:BH:89:LYS:O	30:BH:90:LEU:HD12	2.21	0.40
31:BJ:39:LYS:C	31:BJ:41:LYS:N	2.74	0.40
33:BL:106:GLU:HG2	33:BL:107:PHE:N	2.36	0.40
23:BB:911:A:N6	34:BM:13:HIS:HB3	2.36	0.40
34:BM:70:ASP:C	34:BM:71:LYS:HG3	2.41	0.40
34:BM:82:MET:O	34:BM:84:LYS:HG3	2.21	0.40
34:BM:96:ILE:O	34:BM:97:GLN:HG3	2.21	0.40
35:BN:110:MET:O	35:BN:111:ALA:HB3	2.20	0.40
35:BN:28:LEU:HD22	35:BN:48:VAL:HG11	2.03	0.40
37:BP:2:ASN:ND2	37:BP:2:ASN:N	2.61	0.40
40:BS:24:ILE:O	40:BS:25:ARG:HB2	2.20	0.40
42:BU:69:VAL:CG1	42:BU:77:GLY:H	2.34	0.40
44:BX:36:GLN:OE1	44:BX:37:LEU:HG	2.22	0.40
46:BZ:49:ARG:C	46:BZ:51:VAL:N	2.72	0.40
1:CA:1271:A:H2'	1:CA:1272:G:H8	1.86	0.40
1:CA:1526:G:H2'	1:CA:1527:U:H6	1.85	0.40
1:CA:1530:G:H2'	1:CA:1531:A:C8	2.55	0.40
1:CA:401:C:H2'	1:CA:402:G:H8	1.86	0.40
1:CA:442:G:H2'	1:CA:443:C:H6	1.84	0.40
1:CA:60:A:HO2'	1:CA:61:G:C5'	2.35	0.40
1:CA:651:C:O2'	1:CA:652:U:H5'	2.21	0.40
1:CA:656:G:O2'	1:CA:657:U:H5'	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:977:A:H3'	1:CA:1362:A:N6	2.32	0.40
20:CB:26:MET:SD	20:CB:192:PRO:HD3	2.61	0.40
20:CB:207:ARG:O	20:CB:208:ALA:C	2.59	0.40
20:CB:77:GLU:O	20:CB:80:LYS:HG2	2.21	0.40
9:CJ:57:VAL:HG13	9:CJ:58:ASN:N	2.36	0.40
10:CK:67:GLU:HG3	10:CK:68:ARG:N	2.36	0.40
11:CL:56:LEU:HD12	11:CL:60:PHE:CB	2.50	0.40
15:CP:43:ALA:C	15:CP:45:GLU:H	2.23	0.40
18:CS:9:PHE:O	18:CS:10:ILE:HG23	2.21	0.40
19:CT:38:ILE:HD11	19:CT:82:ILE:CG2	2.47	0.40
48:D1:4:ILE:HG13	48:D1:4:ILE:O	2.21	0.40
22:DA:49:C:H5''	36:DO:101:GLY:HA3	2.03	0.40
23:DB:1019:U:O2'	23:DB:1021:A:C2	2.70	0.40
23:DB:1092:C:C2'	23:DB:1093:G:H5'	2.51	0.40
23:DB:1549:A:H2'	23:DB:1550:C:H6	1.84	0.40
23:DB:1310:G:H1'	23:DB:1611:C:H5'	2.03	0.40
23:DB:1764:C:H2'	23:DB:1765:U:C6	2.56	0.40
23:DB:1792:G:O2'	23:DB:1793:C:H5'	2.21	0.40
23:DB:1864:U:OP1	23:DB:2411:A:H5'	2.21	0.40
23:DB:189:G:H2'	23:DB:205:G:N2	2.37	0.40
23:DB:1936:A:C2	23:DB:1943:U:H5	2.39	0.40
23:DB:2215:C:H2'	23:DB:2216:G:C8	2.56	0.40
23:DB:2221:G:H2'	23:DB:2222:C:H6	1.85	0.40
23:DB:2389:G:H5''	23:DB:2390:U:O4'	2.20	0.40
23:DB:182:A:H1'	23:DB:434:U:H5'	2.03	0.40
23:DB:848:C:H2'	23:DB:849:A:C8	2.56	0.40
22:DA:99:A:N1	23:DB:863:A:C4'	2.85	0.40
25:DC:122:ALA:O	25:DC:124:LYS:HG2	2.21	0.40
25:DC:68:ARG:HH21	25:DC:190:THR:HG23	1.83	0.40
25:DC:9:SER:OG	25:DC:202:ARG:NH2	2.55	0.40
26:DD:79:LEU:CG	26:DD:80:TRP:N	2.83	0.40
31:DJ:133:ALA:C	31:DJ:135:GLN:H	2.23	0.40
31:DJ:33:ALA:O	31:DJ:36:LEU:HB2	2.20	0.40
32:DK:43:ILE:CD1	32:DK:58:LEU:HD21	2.50	0.40
23:DB:2405:G:C5'	33:DL:70:LYS:HE3	2.51	0.40
36:DO:15:ARG:HH12	43:DW:76:ARG:HD2	1.86	0.40
37:DP:50:ARG:HD2	37:DP:99:LEU:HB3	2.03	0.40
39:DR:39:LEU:HG	39:DR:62:GLU:HG2	2.02	0.40
39:DR:73:LYS:HD2	39:DR:73:LYS:H	1.85	0.40
39:DR:76:LYS:CE	39:DR:90:ARG:HD3	2.52	0.40
41:DT:12:ARG:HG2	41:DT:13:ALA:H	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DU:43:LYS:NZ	42:DU:43:LYS:CB	2.84	0.40
42:DU:44:HIS:O	42:DU:46:LYS:HD2	2.21	0.40
42:DU:96:LYS:HB2	42:DU:96:LYS:HE2	1.94	0.40
24:DV:65:VAL:O	24:DV:67:GLY:N	2.51	0.40
23:DB:2279:G:N7	43:DW:10:ARG:HG3	2.36	0.40
43:DW:24:ARG:HE	43:DW:58:LEU:HB2	1.86	0.40
43:DW:48:ALA:HA	43:DW:54:ARG:N	2.35	0.40
46:DZ:1:MET:C	46:DZ:2:LYS:HG3	2.41	0.40
1:AA:123:U:H5''	1:AA:311:C:HO2'	1.86	0.40
1:AA:1343:G:H1'	8:AI:122:ARG:NH1	2.35	0.40
1:AA:175:C:H2'	1:AA:176:C:H6	1.85	0.40
1:AA:239:U:C6	1:AA:239:U:C5'	2.97	0.40
1:AA:471:U:O2'	1:AA:472:U:H5'	2.21	0.40
1:AA:669:G:O2'	1:AA:670:G:H5'	2.22	0.40
2:AC:90:VAL:C	2:AC:93:ILE:HG22	2.42	0.40
3:AD:153:ARG:HG2	3:AD:153:ARG:NH1	2.36	0.40
3:AD:187:ARG:HG3	3:AD:187:ARG:HH11	1.85	0.40
5:AF:88:MET:HE1	5:AF:89:VAL:C	2.41	0.40
6:AG:132:THR:HG23	6:AG:135:LYS:NZ	2.35	0.40
7:AH:26:MET:CB	7:AH:27:PRO:HD2	2.49	0.40
7:AH:76:ARG:HD3	7:AH:79:ARG:HB3	2.02	0.40
11:AL:31:GLY:O	11:AL:78:VAL:HG13	2.22	0.40
12:AM:106:ARG:HD3	12:AM:109:LYS:HD2	2.03	0.40
14:AO:82:GLU:OE1	14:AO:82:GLU:HA	2.21	0.40
16:AQ:10:ARG:NE	16:AQ:56:ASP:O	2.54	0.40
16:AQ:4:ILE:HD12	16:AQ:5:ARG:H	1.85	0.40
18:AS:18:VAL:HG13	18:AS:19:GLU:N	2.36	0.40
48:B1:35:LEU:O	48:B1:35:LEU:HG	2.21	0.40
49:B2:8:SER:OG	49:B2:11:LYS:HB2	2.21	0.40
22:BA:30:C:H2'	22:BA:31:C:H5'	2.03	0.40
22:BA:81:G:H2'	22:BA:82:U:H6	1.85	0.40
23:BB:1017:G:H2'	23:BB:1018:U:H6	1.86	0.40
23:BB:102:U:H4'	23:BB:103:A:OP2	2.13	0.40
23:BB:1185:G:H5''	23:BB:1186:G:OP1	2.21	0.40
23:BB:123:G:O2'	23:BB:124:G:H5'	2.21	0.40
23:BB:135:U:O5'	23:BB:135:U:H6	2.04	0.40
23:BB:1411:U:O2'	23:BB:1412:U:H5'	2.21	0.40
23:BB:154:U:O2'	23:BB:155:A:H5'	2.22	0.40
23:BB:1719:G:O2'	23:BB:1720:U:H5'	2.22	0.40
23:BB:2311:A:H1'	28:BF:76:PHE:HE1	1.87	0.40
23:BB:2313:C:H2'	23:BB:2314:A:H8	1.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2732:G:H3'	23:BB:2733:A:H5'	2.02	0.40
23:BB:2799:A:H4'	23:BB:2800:A:C8	2.56	0.40
23:BB:2838:G:C4	23:BB:2839:G:C8	3.10	0.40
23:BB:428:A:H2'	23:BB:429:A:O4'	2.21	0.40
23:BB:622:G:H2'	23:BB:623:C:C6	2.56	0.40
23:BB:975:A:N6	23:BB:989:G:H1'	2.37	0.40
23:BB:977:G:H5'	23:BB:1155:A:H4'	2.03	0.40
25:BC:175:LEU:HG	25:BC:175:LEU:H	1.74	0.40
25:BC:216:ARG:O	25:BC:218:THR:N	2.53	0.40
25:BC:220:ARG:NH1	25:BC:220:ARG:HG3	2.36	0.40
25:BC:237:ARG:HE	25:BC:237:ARG:C	2.25	0.40
25:BC:259:ASN:O	25:BC:261:ARG:HG2	2.21	0.40
26:BD:103:ASP:C	26:BD:105:LYS:H	2.23	0.40
26:BD:187:LEU:HD23	26:BD:187:LEU:C	2.41	0.40
26:BD:192:ALA:O	26:BD:193:VAL:HG12	2.22	0.40
28:BF:109:ARG:CG	28:BF:136:ILE:HG13	2.50	0.40
28:BF:13:LYS:HE2	28:BF:24:VAL:HG12	2.03	0.40
28:BF:12:VAL:CG1	28:BF:27:VAL:HG21	2.49	0.40
29:BG:131:VAL:HG13	29:BG:133:LYS:CE	2.49	0.40
29:BG:26:LYS:HB3	29:BG:26:LYS:HE2	1.85	0.40
30:BH:28:ASN:O	30:BH:29:PHE:HB2	2.21	0.40
31:BJ:23:LYS:N	31:BJ:23:LYS:HD2	2.33	0.40
31:BJ:60:ASP:OD1	31:BJ:97:PRO:HG2	2.21	0.40
33:BL:65:GLY:O	33:BL:66:PHE:HB3	2.20	0.40
33:BL:94:THR:CG2	33:BL:95:LEU:HD12	2.37	0.40
34:BM:21:ALA:C	34:BM:96:ILE:HG13	2.42	0.40
34:BM:81:ARG:HG3	34:BM:81:ARG:O	2.21	0.40
35:BN:73:ASN:N	35:BN:73:ASN:ND2	2.68	0.40
32:BK:76:VAL:O	37:BP:74:GLN:HG2	2.21	0.40
37:BP:92:ARG:NH2	37:BP:110:LYS:HB3	2.36	0.40
38:BQ:5:ARG:H	38:BQ:5:ARG:HD3	1.86	0.40
38:BQ:92:LYS:HG2	38:BQ:93:ILE:N	2.10	0.40
40:BS:86:MET:HB3	40:BS:94:ASP:HB3	2.03	0.40
42:BU:10:VAL:O	42:BU:11:ILE:HD13	2.22	0.40
24:BV:9:ARG:HH22	24:BV:27:PRO:HB3	1.86	0.40
45:BY:6:ILE:O	45:BY:6:ILE:HG12	2.20	0.40
1:CA:1011:C:H2'	1:CA:1012:A:H8	1.85	0.40
1:CA:119:A:H4'	1:CA:120:A:O4'	2.21	0.40
1:CA:1209:C:O2'	1:CA:1210:C:H5'	2.21	0.40
1:CA:123:U:H2'	1:CA:124:C:H6	1.86	0.40
1:CA:1446:A:N6	1:CA:1447:A:N6	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:144:G:H2'	1:CA:145:G:C8	2.56	0.40
1:CA:213:G:C8	1:CA:214:C:C5	3.09	0.40
1:CA:374:A:OP1	1:CA:452:A:N1	2.55	0.40
1:CA:666:G:H5'	1:CA:726:C:H1'	2.03	0.40
1:CA:99:C:H2'	55:CA:1943:HOH:O	2.21	0.40
3:CD:12:ARG:C	3:CD:37:PRO:HG3	2.42	0.40
3:CD:54:LEU:HD13	3:CD:54:LEU:C	2.41	0.40
4:CE:104:ILE:O	4:CE:104:ILE:HD13	2.21	0.40
4:CE:146:MET:HG2	4:CE:147:ASN:N	2.36	0.40
6:CG:132:THR:O	6:CG:135:LYS:HB3	2.21	0.40
7:CH:54:THR:HG23	7:CH:55:LYS:HE2	2.02	0.40
13:CN:68:ARG:HH12	13:CN:71:GLY:H	1.67	0.40
16:CQ:56:ASP:OD1	16:CQ:80:LYS:HA	2.21	0.40
18:CS:39:ILE:HA	18:CS:43:MET:SD	2.60	0.40
51:D4:28:SER:O	51:D4:29:ALA:HB3	2.21	0.40
23:DB:1228:G:O2'	23:DB:1229:C:H5'	2.21	0.40
23:DB:1266:G:OP2	47:D0:12:ARG:CZ	2.69	0.40
23:DB:1708:C:H2'	23:DB:1709:U:C6	2.57	0.40
23:DB:1723:G:N7	23:DB:1724:G:N7	2.69	0.40
23:DB:1856:U:C2'	23:DB:1857:G:H5'	2.52	0.40
23:DB:1945:G:H2'	23:DB:1946:U:H6	1.82	0.40
23:DB:2186:G:H2'	23:DB:2187:U:O4'	2.21	0.40
23:DB:2673:G:H2'	23:DB:2674:G:H8	1.86	0.40
23:DB:2820:A:O3'	35:DN:2:ARG:NH2	2.53	0.40
23:DB:311:A:HO2'	23:DB:312:G:C5'	2.35	0.40
23:DB:553:G:C2'	23:DB:554:U:H5'	2.52	0.40
25:DC:160:TYR:CD2	25:DC:193:GLU:HG2	2.56	0.40
26:DD:60:VAL:CG2	26:DD:63:PRO:HG2	2.43	0.40
28:DF:157:THR:HG22	28:DF:159:ALA:H	1.87	0.40
31:DJ:23:LYS:HB3	31:DJ:23:LYS:HE2	1.87	0.40
33:DL:45:GLY:C	33:DL:47:ARG:N	2.74	0.40
34:DM:102:LEU:N	34:DM:102:LEU:HD22	2.35	0.40
34:DM:42:THR:O	34:DM:45:GLN:HB2	2.22	0.40
34:DM:5:LYS:O	34:DM:7:THR:HG23	2.21	0.40
34:DM:71:LYS:HG2	34:DM:71:LYS:H	1.58	0.40
36:DO:109:ALA:HA	36:DO:112:GLU:CG	2.51	0.40
37:DP:36:LYS:CG	37:DP:37:LYS:H	2.33	0.40
38:DQ:101:ASP:O	38:DQ:105:PHE:HB2	2.21	0.40
38:DQ:64:ILE:C	38:DQ:66:ALA:N	2.75	0.40
40:DS:64:ALA:HA	40:DS:110:ARG:NH2	2.36	0.40
41:DT:15:HIS:HB3	41:DT:31:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DT:7:LEU:CD2	41:DT:7:LEU:H	2.33	0.40
43:DW:67:LYS:HG2	43:DW:71:LYS:N	2.36	0.40
43:DW:36:ILE:HG12	43:DW:69:GLU:OE2	2.21	0.40
44:DX:26:PHE:N	44:DX:26:PHE:CD1	2.89	0.40
44:DX:41:HIS:CE1	44:DX:43:LEU:HD23	2.56	0.40
46:DZ:49:ARG:HB2	46:DZ:49:ARG:NH1	2.36	0.40
1:AA:1080:A:H2'	1:AA:1081:A:H5'	2.03	0.40
1:AA:1237:C:C4'	1:AA:1334:G:N2	2.85	0.40
1:AA:1340:A:H2'	1:AA:1341:U:C6	2.57	0.40
1:AA:258:G:N2	1:AA:259:G:H1'	2.37	0.40
1:AA:619:U:O2	3:AD:129:VAL:HA	2.22	0.40
1:AA:672:U:H2'	1:AA:673:A:H8	1.83	0.40
1:AA:80:A:C2	1:AA:81:A:H1'	2.57	0.40
1:AA:828:U:H2'	1:AA:829:G:O5'	2.21	0.40
1:AA:82:G:H3'	1:AA:83:C:H4'	2.02	0.40
1:AA:977:A:N1	1:AA:1224:U:O5'	2.54	0.40
1:AA:978:A:H5'	1:AA:1362:A:H62	1.83	0.40
20:AB:98:GLY:HA2	20:AB:101:THR:CG2	2.51	0.40
2:AC:205:GLU:HB2	2:AC:206:ILE:H	1.59	0.40
7:AH:30:LYS:NZ	7:AH:30:LYS:CA	2.78	0.40
11:AL:19:ASN:HB2	11:AL:93:ARG:CZ	2.52	0.40
18:AS:4:LEU:N	18:AS:5:LYS:HE3	2.35	0.40
48:B1:51:ALA:CB	48:B1:52:LYS:HD2	2.52	0.40
50:B3:22:LYS:HA	50:B3:22:LYS:HD3	1.88	0.40
23:BB:1036:G:O2'	23:BB:1037:G:H5'	2.21	0.40
23:BB:1050:A:C2	23:BB:1051:G:H1'	2.57	0.40
23:BB:1254:A:OP2	23:BB:1256:G:H8	2.04	0.40
23:BB:1306:C:O2'	23:BB:1307:A:H5'	2.22	0.40
23:BB:143:C:C2'	23:BB:144:A:C8	3.04	0.40
23:BB:1541:C:H2'	23:BB:1542:U:H6	1.85	0.40
23:BB:1695:G:N7	25:BC:14:HIS:HD2	2.19	0.40
23:BB:1788:C:O2'	23:BB:1789:A:H5'	2.21	0.40
23:BB:20:C:O2'	23:BB:21:A:H5'	2.21	0.40
23:BB:2528:U:P	51:B4:31:PRO:HG2	2.61	0.40
23:BB:1758:U:N3	23:BB:2696:U:H5'	2.35	0.40
23:BB:2887:A:O2'	23:BB:2888:C:H5'	2.21	0.40
23:BB:442:G:N2	23:BB:444:C:C2	2.90	0.40
23:BB:515:A:H3'	23:BB:516:C:C6	2.56	0.40
23:BB:878:A:OP1	23:BB:899:A:N6	2.53	0.40
25:BC:228:ASP:N	25:BC:228:ASP:OD1	2.53	0.40
25:BC:50:THR:C	25:BC:51:ARG:HG3	2.42	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:BD:128:ARG:NH2	26:BD:130:GLN:CD	2.74	0.40
23:BB:2574:G:O2'	26:BD:148:GLN:NE2	2.55	0.40
26:BD:33:ARG:HG3	26:BD:34:VAL:N	2.31	0.40
27:BE:176:ASP:CG	27:BE:177:PRO:HD2	2.41	0.40
27:BE:52:VAL:HG22	27:BE:53:THR:N	2.29	0.40
27:BE:37:ALA:O	27:BE:93:SER:HA	2.21	0.40
30:BH:82:SER:O	30:BH:90:LEU:HD23	2.21	0.40
31:BJ:73:VAL:HG21	31:BJ:75:TYR:OH	2.21	0.40
33:BL:123:ARG:O	33:BL:126:ARG:HG3	2.21	0.40
33:BL:69:ARG:HG2	33:BL:69:ARG:HH11	1.86	0.40
34:BM:33:LEU:H	34:BM:101:VAL:HG22	1.81	0.40
34:BM:71:LYS:O	34:BM:73:ILE:HD12	2.22	0.40
35:BN:65:LEU:HA	35:BN:65:LEU:HD12	1.93	0.40
36:BO:34:HIS:N	36:BO:34:HIS:ND1	2.69	0.40
37:BP:51:ASN:ND2	37:BP:97:TYR:CE2	2.90	0.40
37:BP:67:GLU:O	37:BP:67:GLU:HG3	2.20	0.40
38:BQ:93:ILE:HG23	38:BQ:94:LEU:N	2.31	0.40
38:BQ:94:LEU:HB3	39:BR:13:ARG:HG2	2.03	0.40
40:BS:6:LYS:HZ1	40:BS:104:THR:HG22	1.86	0.40
40:BS:3:THR:HG23	40:BS:107:VAL:O	2.22	0.40
41:BT:24:MET:HG3	41:BT:28:ASN:HA	2.02	0.40
41:BT:68:LYS:HA	41:BT:75:GLY:HA3	2.03	0.40
42:BU:43:LYS:HE2	42:BU:45:GLN:HG2	2.03	0.40
44:BX:14:LEU:HD21	44:BX:57:LEU:HD22	2.02	0.40
41:BT:12:ARG:NH2	44:BX:29:ARG:NH1	2.59	0.40
1:CA:1046:A:H2'	1:CA:1047:G:O4'	2.21	0.40
1:CA:1074:G:H2'	1:CA:1075:U:C6	2.53	0.40
1:CA:1099:G:H2'	1:CA:1100:C:O4'	2.21	0.40
1:CA:1125:U:O2	1:CA:1126:U:H6	2.04	0.40
1:CA:1321:U:H2'	1:CA:1322:C:C5	2.56	0.40
1:CA:1326:U:H2'	1:CA:1327:C:H6	1.86	0.40
1:CA:412:A:H1'	1:CA:413:G:C8	2.45	0.40
1:CA:471:U:O2'	1:CA:472:U:H5'	2.22	0.40
1:CA:598:U:C2	1:CA:599:C:C5	3.09	0.40
1:CA:612:C:H2'	1:CA:613:C:C6	2.56	0.40
1:CA:846:G:H5'	1:CA:847:G:OP2	2.21	0.40
2:CC:8:GLY:C	2:CC:11:LEU:HG	2.41	0.40
2:CC:85:LYS:O	2:CC:89:VAL:HG13	2.21	0.40
1:CA:619:U:O2	3:CD:129:VAL:HG13	2.22	0.40
3:CD:167:PRO:CG	3:CD:170:LEU:HD11	2.35	0.40
3:CD:44:LYS:HZ1	3:CD:46:ARG:HA	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CF:94:HIS:O	5:CF:95:ALA:C	2.59	0.40
7:CH:7:ALA:O	7:CH:11:THR:HG23	2.21	0.40
8:CI:85:ALA:C	8:CI:87:MET:N	2.74	0.40
11:CL:24:GLU:HB2	11:CL:26:CYS:SG	2.61	0.40
11:CL:62:VAL:HG11	11:CL:94:TYR:HE1	1.87	0.40
1:CA:552:U:H5'	11:CL:82:ARG:HH11	1.87	0.40
12:CM:21:ILE:O	12:CM:23:GLY:N	2.55	0.40
12:CM:13:HIS:HA	12:CM:43:LYS:HA	2.02	0.40
13:CN:9:GLU:O	13:CN:13:VAL:HG23	2.20	0.40
14:CO:16:ARG:CZ	14:CO:76:ARG:NH2	2.85	0.40
18:CS:35:ARG:NH2	18:CS:74:ALA:HB3	2.37	0.40
18:CS:39:ILE:HG12	18:CS:68:HIS:O	2.20	0.40
21:CU:3:ILE:HD11	21:CU:19:LYS:HD3	2.04	0.40
47:D0:41:HIS:HB3	47:D0:46:GLY:CA	2.51	0.40
48:D1:33:LEU:N	48:D1:33:LEU:HD12	2.37	0.40
23:DB:242:G:H5'	50:D3:61:LEU:HD23	2.04	0.40
50:D3:4:LYS:O	50:D3:6:VAL:N	2.54	0.40
22:DA:40:U:C2	22:DA:43:C:H5''	2.56	0.40
23:DB:1299:G:C5'	23:DB:1301:A:H1'	2.51	0.40
23:DB:137:U:C5	23:DB:138:U:C4	3.09	0.40
23:DB:1506:U:H2'	23:DB:1507:C:H6	1.87	0.40
23:DB:1989:G:H2'	23:DB:1990:C:O4'	2.21	0.40
23:DB:2143:C:H3'	23:DB:2144:G:C8	2.48	0.40
23:DB:2150:C:H2'	23:DB:2151:U:H6	1.85	0.40
23:DB:2786:U:O2	26:DD:62:LYS:HB3	2.22	0.40
23:DB:622:G:H2'	23:DB:623:C:C6	2.56	0.40
23:DB:689:A:H2'	23:DB:690:G:H8	1.85	0.40
23:DB:876:C:H3'	23:DB:877:A:C4'	2.52	0.40
23:DB:7:G:H2'	23:DB:8:C:O4'	2.21	0.40
25:DC:131:MET:HE2	25:DC:173:LEU:HD11	2.03	0.40
25:DC:22:GLU:HB2	25:DC:202:ARG:NE	2.37	0.40
25:DC:51:ARG:O	25:DC:52:HIS:O	2.39	0.40
26:DD:175:LEU:HB3	26:DD:189:VAL:HG12	2.02	0.40
26:DD:72:GLY:O	26:DD:91:THR:OG1	2.38	0.40
26:DD:81:GLU:O	26:DD:82:PHE:CB	2.69	0.40
26:DD:87:GLY:HA3	26:DD:89:GLU:OE1	2.21	0.40
31:DJ:124:VAL:O	31:DJ:125:TYR:HB2	2.20	0.40
31:DJ:15:TRP:HB3	31:DJ:138:GLN:O	2.22	0.40
34:DM:81:ARG:HB2	34:DM:81:ARG:HH11	1.86	0.40
35:DN:81:ASN:O	35:DN:85:PRO:HG3	2.22	0.40
36:DO:39:VAL:O	36:DO:39:VAL:HG22	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:DP:38:ARG:HA	37:DP:38:ARG:HD2	1.61	0.40
38:DQ:30:VAL:O	38:DQ:31:TYR:HB2	2.21	0.40
38:DQ:71:ASN:O	38:DQ:73:ILE:N	2.54	0.40
38:DQ:86:SER:C	38:DQ:88:GLU:H	2.24	0.40
41:DT:87:LEU:HD13	41:DT:93:LEU:HD13	2.02	0.40
43:DW:77:LYS:C	43:DW:79:ILE:H	2.24	0.40
1:AA:1021:A:H2'	1:AA:1022:A:C8	2.56	0.40
1:AA:1057:G:H2'	1:AA:1058:G:O4'	2.21	0.40
1:AA:1095:U:O2'	1:AA:1096:C:H5'	2.22	0.40
1:AA:1238:A:H2	1:AA:1241:G:H1'	1.85	0.40
1:AA:123:U:H2'	1:AA:124:C:H6	1.86	0.40
1:AA:1371:G:O3'	8:AI:70:GLY:HA3	2.20	0.40
1:AA:1458:G:O3'	19:AT:22:SER:HA	2.22	0.40
1:AA:201:G:O2'	1:AA:202:G:H5'	2.21	0.40
1:AA:438:U:H4'	3:AD:119:HIS:HD2	1.85	0.40
1:AA:454:G:H2'	1:AA:455:G:H8	1.86	0.40
2:AC:154:GLY:CA	2:AC:163:ARG:H	2.33	0.40
2:AC:5:HIS:NE2	2:AC:7:ASN:HB2	2.37	0.40
3:AD:84:ASN:O	3:AD:88:ASN:ND2	2.54	0.40
4:AE:95:MET:HE2	4:AE:114:LEU:HD11	2.03	0.40
8:AI:16:ALA:HB1	8:AI:78:ILE:HA	2.03	0.40
1:AA:973:G:O2'	9:AJ:56:HIS:HA	2.21	0.40
14:AO:24:THR:O	14:AO:28:VAL:HG23	2.21	0.40
18:AS:30:LEU:N	18:AS:30:LEU:HD22	2.36	0.40
19:AT:17:ARG:HD2	19:AT:17:ARG:C	2.42	0.40
48:B1:10:LEU:HD11	48:B1:52:LYS:HD3	2.04	0.40
49:B2:42:LEU:HG	49:B2:43:THR:H	1.86	0.40
49:B2:12:ARG:NH1	49:B2:44:VAL:O	2.55	0.40
23:BB:2363:G:P	50:B3:39:ARG:HD3	2.62	0.40
51:B4:23:ILE:H	51:B4:23:ILE:HG12	1.64	0.40
23:BB:1010:A:H4'	23:BB:1152:C:O2'	2.22	0.40
23:BB:1018:U:H2'	23:BB:1019:U:O4'	2.22	0.40
23:BB:116:C:H2'	23:BB:117:G:C8	2.57	0.40
23:BB:1708:C:H2'	23:BB:1709:U:C6	2.56	0.40
23:BB:1893:C:H2'	23:BB:1894:C:O4'	2.21	0.40
23:BB:18:U:O2'	23:BB:19:A:H5'	2.21	0.40
23:BB:2080:A:H2'	23:BB:2081:U:H6	1.86	0.40
23:BB:2155:U:H6	23:BB:2155:U:O5'	2.05	0.40
23:BB:2179:C:H2'	23:BB:2180:U:C6	2.56	0.40
23:BB:2312:U:H2'	23:BB:2313:C:H5'	2.03	0.40
23:BB:2377:A:H2'	23:BB:2378:A:C8	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:BB:2444:G:OP1	27:BE:62:GLN:NE2	2.55	0.40
23:BB:27:G:O2'	23:BB:513:A:N6	2.54	0.40
23:BB:2822:G:OP1	26:BD:166:GLY:CA	2.69	0.40
23:BB:553:G:H2'	23:BB:554:U:C6	2.56	0.40
23:BB:920:A:H2'	23:BB:921:C:C6	2.55	0.40
27:BE:6:LYS:HZ1	27:BE:119:ILE:N	2.19	0.40
27:BE:48:THR:CA	27:BE:86:ALA:HB1	2.51	0.40
29:BG:35:THR:C	29:BG:36:LEU:HD12	2.42	0.40
29:BG:40:VAL:CG1	29:BG:41:GLU:N	2.84	0.40
30:BH:57:LYS:HA	30:BH:61:VAL:HB	2.02	0.40
52:BI:75:ALA:HB3	52:BI:131:THR:HG21	2.03	0.40
31:BJ:110:PRO:O	31:BJ:111:LYS:HB2	2.22	0.40
33:BL:120:VAL:HG12	33:BL:138:ALA:HB3	2.02	0.40
33:BL:55:MET:SD	33:BL:55:MET:N	2.93	0.40
35:BN:22:ARG:NE	35:BN:69:ARG:HD3	2.37	0.40
35:BN:28:LEU:HG	35:BN:34:ILE:CD1	2.51	0.40
35:BN:41:ALA:C	35:BN:43:GLU:H	2.24	0.40
36:BO:40:ILE:N	36:BO:40:ILE:HD13	2.35	0.40
39:BR:41:ILE:HD12	39:BR:43:ASN:H	1.86	0.40
40:BS:13:SER:OG	40:BS:16:LYS:HD3	2.21	0.40
41:BT:3:ARG:HE	41:BT:8:LEU:CD2	2.34	0.40
42:BU:8:ASP:O	42:BU:10:VAL:HG13	2.21	0.40
42:BU:13:LEU:O	42:BU:15:GLY:N	2.54	0.40
42:BU:25:LYS:CG	42:BU:35:VAL:HA	2.41	0.40
24:BV:35:GLU:HG3	24:BV:93:ARG:NE	2.35	0.40
43:BW:59:PHE:HE2	43:BW:81:ILE:CD1	2.35	0.40
44:BX:19:LEU:C	44:BX:21:LEU:H	2.25	0.40
46:BZ:34:LEU:H	46:BZ:44:PHE:CB	2.34	0.40
1:CA:1336:C:H4'	1:CA:1337:G:O5'	2.21	0.40
1:CA:37:U:O2'	1:CA:38:G:H5'	2.21	0.40
1:CA:546:A:H4'	1:CA:548:G:O3'	2.21	0.40
1:CA:551:U:H2'	1:CA:552:U:H6	1.86	0.40
1:CA:621:A:H2'	1:CA:622:A:H8	1.82	0.40
1:CA:812:G:OP1	1:CA:812:G:C4'	2.70	0.40
20:CB:69:VAL:CG2	20:CB:160:LEU:HD11	2.51	0.40
20:CB:48:MET:CB	20:CB:199:ILE:HG22	2.52	0.40
20:CB:186:VAL:O	20:CB:200:PRO:HA	2.22	0.40
20:CB:64:GLY:O	20:CB:66:ILE:HG12	2.22	0.40
4:CE:61:LYS:HG3	4:CE:62:ALA:N	2.35	0.40
5:CF:89:VAL:HG13	5:CF:89:VAL:O	2.20	0.40
11:CL:20:VAL:O	11:CL:20:VAL:HG23	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CM:22:TYR:O	12:CM:24:VAL:N	2.55	0.40
13:CN:24:ALA:C	13:CN:26:LEU:H	2.25	0.40
13:CN:48:GLN:HG3	13:CN:48:GLN:H	1.61	0.40
14:CO:10:ILE:HG21	14:CO:30:LEU:HD21	2.03	0.40
16:CQ:60:ILE:HB	16:CQ:73:THR:O	2.21	0.40
16:CQ:57:VAL:HG12	16:CQ:78:VAL:HB	2.03	0.40
18:CS:28:LYS:N	18:CS:28:LYS:HD2	2.37	0.40
19:CT:68:LYS:HB2	19:CT:70:LYS:CD	2.51	0.40
48:D1:10:LEU:H	48:D1:24:LYS:CB	2.31	0.40
50:D3:12:ARG:O	50:D3:13:PHE:CB	2.69	0.40
23:DB:1176:U:O5'	23:DB:1176:U:H6	2.04	0.40
23:DB:1275:A:N3	23:DB:1275:A:C3'	2.79	0.40
23:DB:1300:G:H4'	23:DB:1301:A:O5'	2.21	0.40
23:DB:2586:U:H2'	23:DB:2587:A:H8	1.84	0.40
23:DB:769:U:H2'	23:DB:770:G:C8	2.57	0.40
23:DB:988:A:H5''	45:DY:10:ARG:HD3	2.03	0.40
25:DC:28:PRO:CD	25:DC:79:ARG:HE	2.34	0.40
26:DD:18:ASP:O	26:DD:20:VAL:N	2.54	0.40
26:DD:95:SER:HB3	26:DD:96:ILE:H	1.65	0.40
27:DE:51:GLU:HG2	27:DE:52:VAL:HG13	2.03	0.40
28:DF:116:LEU:HD22	28:DF:129:MET:CE	2.52	0.40
29:DG:171:LYS:HZ3	29:DG:173:ALA:HA	1.86	0.40
30:DH:122:LEU:HA	30:DH:146:VAL:CG2	2.49	0.40
31:DJ:7:LYS:CE	31:DJ:45:THR:HG21	2.52	0.40
31:DJ:7:LYS:O	31:DJ:48:VAL:HG21	2.20	0.40
31:DJ:90:GLU:C	31:DJ:92:MET:N	2.73	0.40
32:DK:5:GLN:O	32:DK:20:MET:HG3	2.21	0.40
32:DK:2:ILE:HD12	32:DK:2:ILE:N	2.35	0.40
33:DL:77:ILE:HD11	33:DL:108:ALA:O	2.21	0.40
23:DB:1275:A:C8	35:DN:16:HIS:HB2	2.57	0.40
55:DB:3475:HOH:O	35:DN:4:ARG:HA	2.21	0.40
35:DN:70:THR:HG23	35:DN:72:ASP:HB3	2.04	0.40
22:DA:116:G:H4'	36:DO:54:VAL:HG22	2.02	0.40
37:DP:59:THR:HG23	37:DP:76:HIS:NE2	2.36	0.40
38:DQ:50:ARG:HH22	38:DQ:53:LYS:CE	2.34	0.40
38:DQ:57:ARG:O	38:DQ:61:ILE:HG13	2.21	0.40
38:DQ:86:SER:HB3	38:DQ:88:GLU:OE2	2.21	0.40
39:DR:74:ILE:HD11	39:DR:76:LYS:HE2	2.03	0.40
40:DS:36:LEU:O	40:DS:39:THR:HG22	2.21	0.40
41:DT:19:LYS:HA	41:DT:19:LYS:HD2	1.88	0.40
41:DT:70:HIS:CG	41:DT:71:GLY:H	2.39	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:DX:22:LEU:HD23	44:DX:22:LEU:HA	1.97	0.40
46:DZ:3:LYS:HG2	46:DZ:51:VAL:HG21	2.04	0.40
1:AA:1226:C:O2'	1:AA:1227:A:P	2.79	0.40
1:AA:1315:U:H3'	1:AA:1316:G:C8	2.56	0.40
1:AA:137:U:H2'	1:AA:138:G:C8	2.57	0.40
1:AA:607:A:H2'	1:AA:608:A:H8	1.86	0.40
2:AC:155:ARG:CZ	2:AC:192:TYR:HB2	2.51	0.40
2:AC:10:ARG:O	2:AC:15:LYS:HB2	2.20	0.40
1:AA:620:C:H1'	3:AD:131:ILE:HG21	2.03	0.40
3:AD:25:ARG:HD3	3:AD:26:ALA:N	2.30	0.40
4:AE:19:ARG:HA	4:AE:32:PHE:HA	2.03	0.40
8:AI:33:SER:C	8:AI:35:GLU:N	2.74	0.40
8:AI:38:PHE:HD2	8:AI:38:PHE:HA	1.77	0.40
8:AI:57:VAL:C	8:AI:59:LYS:H	2.25	0.40
9:AJ:92:LEU:H	9:AJ:92:LEU:CD2	2.29	0.40
10:AK:35:ASP:HB3	10:AK:41:LEU:HD11	2.04	0.40
11:AL:66:ILE:CD1	11:AL:73:LEU:HD12	2.51	0.40
17:AR:46:THR:O	17:AR:50:TYR:HB2	2.21	0.40
47:B0:9:ARG:O	47:B0:10:SER:HB3	2.22	0.40
47:B0:35:GLU:CG	47:B0:36:LYS:H	2.34	0.40
48:B1:44:GLN:OE1	48:B1:46:VAL:HB	2.22	0.40
48:B1:8:ILE:HB	48:B1:9:LYS:H	1.60	0.40
23:BB:1345:C:H5'	23:BB:1396:U:H5	1.86	0.40
23:BB:1655:A:H2'	23:BB:1656:C:O4'	2.22	0.40
23:BB:1723:G:N7	23:BB:1724:G:N7	2.70	0.40
23:BB:1736:U:H2'	23:BB:1737:G:O4'	2.20	0.40
23:BB:1801:A:H5'	23:BB:2203:U:O2'	2.21	0.40
23:BB:2331:G:H2'	23:BB:2332:C:C6	2.57	0.40
23:BB:2617:U:C4	23:BB:2618:G:N7	2.90	0.40
23:BB:539:G:O2'	23:BB:540:C:H5'	2.22	0.40
23:BB:711:G:C2'	23:BB:712:G:H5'	2.51	0.40
26:BD:168:GLU:HG2	26:BD:169:ARG:N	2.37	0.40
26:BD:175:LEU:HB3	26:BD:189:VAL:CG1	2.52	0.40
26:BD:26:VAL:O	26:BD:27:ILE:C	2.60	0.40
26:BD:46:ARG:HG3	26:BD:47:ALA:N	2.36	0.40
27:BE:193:VAL:HG11	27:BE:201:ALA:HA	2.02	0.40
30:BH:19:VAL:O	30:BH:20:ASN:C	2.59	0.40
30:BH:32:PRO:C	30:BH:34:GLY:H	2.24	0.40
52:BI:4:VAL:O	52:BI:4:VAL:HG13	2.21	0.40
31:BJ:118:MET:O	31:BJ:119:PHE:C	2.59	0.40
31:BJ:25:LEU:HG	31:BJ:64:VAL:N	2.25	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:BL:93:ASN:O	33:BL:94:THR:HB	2.21	0.40
34:BM:126:ILE:HG13	34:BM:127:LYS:N	2.25	0.40
35:BN:116:VAL:HG22	35:BN:117:ASP:N	2.27	0.40
36:BO:17:LYS:HG2	43:BW:77:LYS:HD2	2.02	0.40
37:BP:89:GLY:O	37:BP:90:ALA:C	2.60	0.40
38:BQ:68:ALA:HB2	38:BQ:105:PHE:HZ	1.87	0.40
41:BT:2:ILE:O	41:BT:5:GLU:OE1	2.39	0.40
24:BV:82:TYR:HB3	34:BM:36:VAL:HG12	2.04	0.40
46:BZ:48:GLN:O	46:BZ:51:VAL:HB	2.20	0.40
46:BZ:6:HIS:H	46:BZ:51:VAL:HG13	1.86	0.40
1:CA:1284:C:H3'	1:CA:1285:A:H8	1.86	0.40
1:CA:1475:G:H4'	23:DB:1689:A:H4'	2.03	0.40
1:CA:1515:G:O2'	1:CA:1516:G:H5'	2.21	0.40
1:CA:1503:A:C8	1:CA:1531:A:H1'	2.56	0.40
1:CA:258:G:N2	1:CA:259:G:H1'	2.37	0.40
1:CA:279:A:C5'	1:CA:280:C:H3'	2.48	0.40
1:CA:415:A:N1	1:CA:428:G:O6	2.54	0.40
2:CC:127:VAL:CG2	2:CC:128:MET:H	2.25	0.40
3:CD:105:GLY:HA3	3:CD:158:LEU:HD23	2.03	0.40
3:CD:58:GLN:CA	3:CD:58:GLN:HE21	2.16	0.40
5:CF:11:HIS:HB3	5:CF:14:GLN:HB2	2.02	0.40
5:CF:59:TYR:HD1	5:CF:59:TYR:H	1.68	0.40
7:CH:55:LYS:CA	7:CH:55:LYS:HZ2	2.24	0.40
8:CI:99:LYS:HG2	8:CI:99:LYS:O	2.21	0.40
11:CL:53:ARG:HG2	11:CL:63:THR:HG22	2.03	0.40
1:CA:525:C:OP1	11:CL:87:LYS:HE2	2.21	0.40
12:CM:56:ARG:HA	12:CM:59:VAL:CG1	2.51	0.40
16:CQ:11:VAL:HG13	16:CQ:20:ILE:CG2	2.47	0.40
1:CA:265:G:H5'	16:CQ:65:PRO:O	2.20	0.40
16:CQ:75:VAL:CG2	16:CQ:76:ARG:N	2.85	0.40
18:CS:3:SER:C	18:CS:4:LEU:HD12	2.42	0.40
21:CU:32:ARG:O	21:CU:32:ARG:HG2	2.21	0.40
47:D0:53:VAL:CG1	47:D0:54:ILE:N	2.85	0.40
48:D1:24:LYS:CB	48:D1:24:LYS:HZ2	2.33	0.40
49:D2:4:THR:O	49:D2:5:PHE:HB2	2.21	0.40
51:D4:2:LYS:HB2	51:D4:2:LYS:HZ2	1.82	0.40
51:D4:36:ARG:HD2	51:D4:37:GLN:O	2.22	0.40
22:DA:115:A:H2'	22:DA:116:G:O4'	2.22	0.40
22:DA:23:G:C2	22:DA:24:G:N1	2.90	0.40
22:DA:25:U:O5'	22:DA:26:C:OP1	2.40	0.40
23:DB:1050:A:O2'	23:DB:1051:G:H5'	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:DB:1099:G:N7	52:DI:3:LYS:CD	2.83	0.40
23:DB:1537:G:C2'	23:DB:1538:G:H5'	2.51	0.40
23:DB:1769:U:O2'	23:DB:1770:G:H5'	2.22	0.40
23:DB:2379:G:H2'	23:DB:2380:C:H6	1.87	0.40
23:DB:239:C:H2'	23:DB:240:C:H6	1.87	0.40
23:DB:2617:U:O2'	23:DB:2618:G:H5'	2.22	0.40
23:DB:265:A:H2'	23:DB:266:G:H1'	2.03	0.40
23:DB:2759:G:H2'	23:DB:2760:C:O4'	2.21	0.40
23:DB:2799:A:H4'	23:DB:2800:A:C8	2.56	0.40
23:DB:297:G:H2'	23:DB:298:G:C8	2.56	0.40
23:DB:44:A:H2'	23:DB:45:G:O4'	2.22	0.40
23:DB:562:U:H2'	23:DB:572:A:O4'	2.21	0.40
23:DB:649:G:H2'	23:DB:650:C:C6	2.56	0.40
23:DB:811:U:H3'	33:DL:32:GLY:O	2.21	0.40
25:DC:19:VAL:CG1	25:DC:20:ASN:H	2.33	0.40
25:DC:247:TRP:HZ2	25:DC:254:LYS:HZ3	1.67	0.40
25:DC:50:THR:O	25:DC:51:ARG:C	2.59	0.40
27:DE:138:LEU:HB2	27:DE:143:LEU:CD1	2.51	0.40
27:DE:198:GLU:HB3	27:DE:199:MET:H	1.59	0.40
29:DG:9:VAL:HB	29:DG:11:PRO:HD3	2.03	0.40
30:DH:90:LEU:HD22	30:DH:122:LEU:HB3	2.03	0.40
52:DI:78:LEU:HD13	52:DI:108:ILE:HG23	2.03	0.40
35:DN:2:ARG:CZ	35:DN:4:ARG:NH1	2.85	0.40
36:DO:62:LEU:HD12	36:DO:62:LEU:H	1.87	0.40
32:DK:76:VAL:H	37:DP:74:GLN:HB2	1.86	0.40
40:DS:86:MET:HG3	40:DS:87:PRO:CD	2.51	0.40
41:DT:3:ARG:HB3	41:DT:5:GLU:OE1	2.21	0.40
42:DU:43:LYS:HZ3	42:DU:53:GLN:HE21	1.70	0.40
43:DW:43:LYS:CD	43:DW:76:ARG:HA	2.49	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	204/233 (88%)	144 (71%)	44 (22%)	16 (8%)	1	12
2	CC	204/233 (88%)	138 (68%)	46 (22%)	20 (10%)	1	8
3	AD	203/206 (98%)	133 (66%)	49 (24%)	21 (10%)	0	8
3	CD	203/206 (98%)	137 (68%)	49 (24%)	17 (8%)	1	11
4	AE	148/167 (89%)	110 (74%)	30 (20%)	8 (5%)	2	21
4	CE	148/167 (89%)	109 (74%)	30 (20%)	9 (6%)	2	18
5	AF	98/135 (73%)	68 (69%)	25 (26%)	5 (5%)	2	22
5	CF	98/135 (73%)	72 (74%)	18 (18%)	8 (8%)	1	11
6	AG	148/179 (83%)	105 (71%)	35 (24%)	8 (5%)	2	21
6	CG	150/179 (84%)	101 (67%)	37 (25%)	12 (8%)	1	12
7	AH	127/130 (98%)	108 (85%)	14 (11%)	5 (4%)	3	30
7	CH	127/130 (98%)	91 (72%)	29 (23%)	7 (6%)	2	20
8	AI	125/130 (96%)	84 (67%)	30 (24%)	11 (9%)	1	10
8	CI	125/130 (96%)	82 (66%)	30 (24%)	13 (10%)	0	7
9	AJ	96/103 (93%)	65 (68%)	20 (21%)	11 (12%)	0	6
9	CJ	96/103 (93%)	59 (62%)	18 (19%)	19 (20%)	0	2
10	AK	115/129 (89%)	75 (65%)	32 (28%)	8 (7%)	1	14
10	CK	115/129 (89%)	76 (66%)	30 (26%)	9 (8%)	1	12
11	AL	121/124 (98%)	73 (60%)	35 (29%)	13 (11%)	0	7
11	CL	121/124 (98%)	76 (63%)	28 (23%)	17 (14%)	0	4
12	AM	112/118 (95%)	89 (80%)	13 (12%)	10 (9%)	1	10
12	CM	111/118 (94%)	82 (74%)	17 (15%)	12 (11%)	0	7
13	AN	92/101 (91%)	64 (70%)	19 (21%)	9 (10%)	1	8
13	CN	92/101 (91%)	50 (54%)	26 (28%)	16 (17%)	0	2
14	AO	86/89 (97%)	66 (77%)	19 (22%)	1 (1%)	15	56
14	CO	86/89 (97%)	69 (80%)	15 (17%)	2 (2%)	7	43
15	AP	80/82 (98%)	60 (75%)	14 (18%)	6 (8%)	1	13
15	CP	78/82 (95%)	55 (70%)	14 (18%)	9 (12%)	0	6
16	AQ	78/84 (93%)	49 (63%)	25 (32%)	4 (5%)	2	22
16	CQ	79/84 (94%)	58 (73%)	15 (19%)	6 (8%)	1	12
17	AR	53/75 (71%)	31 (58%)	16 (30%)	6 (11%)	0	6
17	CR	53/75 (71%)	36 (68%)	13 (24%)	4 (8%)	1	13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
18	AS	77/92 (84%)	54 (70%)	12 (16%)	11 (14%)	0	3
18	CS	78/92 (85%)	50 (64%)	17 (22%)	11 (14%)	0	3
19	AT	83/87 (95%)	68 (82%)	12 (14%)	3 (4%)	4	32
19	CT	83/87 (95%)	62 (75%)	16 (19%)	5 (6%)	2	18
20	AB	216/241 (90%)	148 (68%)	52 (24%)	16 (7%)	1	13
20	CB	216/241 (90%)	149 (69%)	41 (19%)	26 (12%)	0	6
21	AU	49/71 (69%)	23 (47%)	12 (24%)	14 (29%)	0	0
21	CU	49/71 (69%)	26 (53%)	18 (37%)	5 (10%)	1	8
24	BV	92/94 (98%)	61 (66%)	24 (26%)	7 (8%)	1	12
24	DV	92/94 (98%)	64 (70%)	23 (25%)	5 (5%)	2	21
25	BC	265/273 (97%)	94 (36%)	95 (36%)	76 (29%)	0	0
25	DC	265/273 (97%)	97 (37%)	101 (38%)	67 (25%)	0	1
26	BD	207/209 (99%)	87 (42%)	68 (33%)	52 (25%)	0	1
26	DD	207/209 (99%)	91 (44%)	72 (35%)	44 (21%)	0	1
27	BE	199/201 (99%)	101 (51%)	57 (29%)	41 (21%)	0	1
27	DE	199/201 (99%)	90 (45%)	63 (32%)	46 (23%)	0	1
28	BF	176/179 (98%)	99 (56%)	42 (24%)	35 (20%)	0	2
28	DF	176/179 (98%)	98 (56%)	49 (28%)	29 (16%)	0	2
29	BG	174/177 (98%)	112 (64%)	41 (24%)	21 (12%)	0	6
29	DG	174/177 (98%)	108 (62%)	51 (29%)	15 (9%)	1	11
30	BH	147/149 (99%)	83 (56%)	49 (33%)	15 (10%)	1	8
30	DH	147/149 (99%)	83 (56%)	46 (31%)	18 (12%)	0	5
31	BJ	138/142 (97%)	68 (49%)	43 (31%)	27 (20%)	0	2
31	DJ	138/142 (97%)	72 (52%)	36 (26%)	30 (22%)	0	1
32	BK	119/123 (97%)	73 (61%)	27 (23%)	19 (16%)	0	2
32	DK	119/123 (97%)	70 (59%)	30 (25%)	19 (16%)	0	2
33	BL	142/144 (99%)	56 (39%)	47 (33%)	39 (28%)	0	0
33	DL	142/144 (99%)	68 (48%)	35 (25%)	39 (28%)	0	0
34	BM	134/136 (98%)	64 (48%)	37 (28%)	33 (25%)	0	1
34	DM	134/136 (98%)	71 (53%)	43 (32%)	20 (15%)	0	3
35	BN	125/127 (98%)	68 (54%)	41 (33%)	16 (13%)	0	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
35	DN	125/127 (98%)	86 (69%)	29 (23%)	10 (8%)	1	12
36	BO	115/117 (98%)	61 (53%)	34 (30%)	20 (17%)	0	2
36	DO	115/117 (98%)	62 (54%)	33 (29%)	20 (17%)	0	2
37	BP	112/115 (97%)	37 (33%)	38 (34%)	37 (33%)	0	0
37	DP	112/115 (97%)	42 (38%)	37 (33%)	33 (30%)	0	0
38	BQ	115/118 (98%)	79 (69%)	23 (20%)	13 (11%)	0	6
38	DQ	115/118 (98%)	78 (68%)	21 (18%)	16 (14%)	0	4
39	BR	101/103 (98%)	39 (39%)	38 (38%)	24 (24%)	0	1
39	DR	101/103 (98%)	42 (42%)	30 (30%)	29 (29%)	0	0
40	BS	108/110 (98%)	58 (54%)	34 (32%)	16 (15%)	0	3
40	DS	108/110 (98%)	62 (57%)	26 (24%)	20 (18%)	0	2
41	BT	97/100 (97%)	38 (39%)	42 (43%)	17 (18%)	0	2
41	DT	97/100 (97%)	44 (45%)	28 (29%)	25 (26%)	0	1
42	BU	100/104 (96%)	36 (36%)	43 (43%)	21 (21%)	0	1
42	DU	100/104 (96%)	47 (47%)	34 (34%)	19 (19%)	0	2
43	BW	82/85 (96%)	35 (43%)	22 (27%)	25 (30%)	0	0
43	DW	82/85 (96%)	28 (34%)	31 (38%)	23 (28%)	0	0
44	BX	61/63 (97%)	20 (33%)	30 (49%)	11 (18%)	0	2
44	DX	61/63 (97%)	37 (61%)	15 (25%)	9 (15%)	0	3
45	BY	56/59 (95%)	29 (52%)	17 (30%)	10 (18%)	0	2
45	DY	56/59 (95%)	38 (68%)	15 (27%)	3 (5%)	2	21
46	BZ	68/70 (97%)	32 (47%)	23 (34%)	13 (19%)	0	2
46	DZ	68/70 (97%)	36 (53%)	22 (32%)	10 (15%)	0	3
47	B0	54/57 (95%)	27 (50%)	17 (32%)	10 (18%)	0	2
47	D0	54/57 (95%)	25 (46%)	20 (37%)	9 (17%)	0	2
48	B1	52/55 (94%)	22 (42%)	20 (38%)	10 (19%)	0	2
48	D1	52/55 (94%)	21 (40%)	18 (35%)	13 (25%)	0	1
49	B2	44/46 (96%)	22 (50%)	15 (34%)	7 (16%)	0	2
49	D2	44/46 (96%)	22 (50%)	10 (23%)	12 (27%)	0	0
50	B3	62/65 (95%)	26 (42%)	28 (45%)	8 (13%)	0	5
50	D3	62/65 (95%)	30 (48%)	21 (34%)	11 (18%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
51	B4	36/38 (95%)	17 (47%)	9 (25%)	10 (28%)	0	0
51	D4	36/38 (95%)	13 (36%)	10 (28%)	13 (36%)	0	0
52	BI	139/142 (98%)	122 (88%)	12 (9%)	5 (4%)	4	32
52	DI	139/142 (98%)	119 (86%)	15 (11%)	5 (4%)	4	32
All	All	11263/11954 (94%)	6605 (59%)	2995 (27%)	1663 (15%)	0	3

All (1663) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	AC	2	GLN
2	AC	81	GLU
2	AC	91	ALA
3	AD	18	LEU
3	AD	25	ARG
3	AD	191	SER
5	AF	92	THR
8	AI	8	THR
9	AJ	57	VAL
10	AK	56	LYS
10	AK	126	ARG
11	AL	23	LEU
11	AL	72	ASN
11	AL	104	SER
12	AM	6	ILE
12	AM	14	ALA
15	AP	79	ASN
15	AP	81	ALA
20	AB	19	THR
20	AB	22	TRP
20	AB	163	ILE
21	AU	14	ALA
21	AU	22	CYS
25	BC	21	PRO
25	BC	22	GLU
25	BC	28	PRO
25	BC	29	PHE
25	BC	31	PRO
25	BC	32	LEU
25	BC	47	ARG
25	BC	63	ILE
25	BC	64	VAL

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Mol	Chain	Res	Type
25	BC	67	LYS
25	BC	68	ARG
25	BC	72	GLY
25	BC	97	ASP
25	BC	103	ILE
25	BC	114	GLN
25	BC	124	LYS
25	BC	135	PRO
25	BC	140	VAL
25	BC	141	HIS
25	BC	149	LYS
25	BC	156	SER
25	BC	176	ARG
25	BC	227	VAL
25	BC	246	PRO
25	BC	266	ILE
26	BD	9	VAL
26	BD	14	ILE
26	BD	33	ARG
26	BD	59	ARG
26	BD	84	LEU
26	BD	114	LYS
26	BD	127	PHE
26	BD	128	ARG
26	BD	130	GLN
26	BD	139	SER
26	BD	140	HIS
26	BD	141	ARG
26	BD	150	GLN
26	BD	154	LYS
26	BD	155	VAL
26	BD	170	VAL
27	BE	4	VAL
27	BE	44	ARG
27	BE	52	VAL
27	BE	59	PRO
27	BE	84	THR
27	BE	117	ARG
27	BE	129	PRO
27	BE	165	HIS
27	BE	196	VAL
28	BF	31	GLU

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Mol	Chain	Res	Type
28	BF	46	LYS
28	BF	66	ILE
28	BF	73	VAL
28	BF	76	PHE
28	BF	131	VAL
28	BF	139	GLU
28	BF	149	ARG
29	BG	6	ALA
29	BG	25	ILE
29	BG	91	VAL
29	BG	151	ARG
30	BH	29	PHE
30	BH	37	VAL
30	BH	38	PRO
30	BH	41	LYS
31	BJ	4	PHE
31	BJ	43	GLU
31	BJ	44	TYR
31	BJ	64	VAL
31	BJ	73	VAL
31	BJ	100	VAL
31	BJ	101	ILE
31	BJ	119	PHE
31	BJ	137	PRO
32	BK	25	LEU
32	BK	35	VAL
32	BK	53	LYS
32	BK	71	ARG
32	BK	72	PRO
32	BK	89	ASN
32	BK	110	GLU
32	BK	120	PRO
33	BL	17	LYS
33	BL	56	PRO
33	BL	62	PRO
33	BL	73	ILE
33	BL	84	LYS
33	BL	85	VAL
33	BL	90	VAL
33	BL	91	ASP
33	BL	92	LEU
33	BL	104	GLN

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Mol	Chain	Res	Type
33	BL	118	THR
33	BL	119	PRO
33	BL	126	ARG
33	BL	127	VAL
33	BL	141	LYS
34	BM	3	GLN
34	BM	4	PRO
34	BM	14	LYS
34	BM	69	PRO
34	BM	80	VAL
34	BM	81	ARG
34	BM	98	PRO
34	BM	109	PRO
35	BN	3	HIS
35	BN	4	ARG
35	BN	8	ARG
35	BN	37	THR
35	BN	89	SER
35	BN	107	ASN
35	BN	113	ILE
35	BN	116	VAL
36	BO	37	ALA
36	BO	42	PRO
36	BO	54	VAL
36	BO	63	LYS
37	BP	23	ASP
37	BP	24	THR
37	BP	49	ILE
37	BP	55	HIS
37	BP	60	VAL
37	BP	82	SER
37	BP	83	ILE
37	BP	85	VAL
37	BP	90	ALA
37	BP	91	VAL
37	BP	94	ALA
37	BP	111	GLU
38	BQ	34	ALA
38	BQ	89	ILE
38	BQ	92	LYS
39	BR	15	SER
39	BR	24	LYS

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Mol	Chain	Res	Type
39	BR	47	VAL
39	BR	59	ILE
39	BR	74	ILE
39	BR	101	ILE
40	BS	6	LYS
40	BS	103	ILE
41	BT	32	LEU
41	BT	63	VAL
42	BU	4	ILE
42	BU	21	ARG
42	BU	63	ALA
42	BU	92	VAL
43	BW	11	ASN
43	BW	18	LYS
43	BW	44	PHE
43	BW	53	GLY
43	BW	66	VAL
43	BW	67	LYS
43	BW	75	ASN
43	BW	81	ILE
44	BX	24	GLU
44	BX	33	ALA
44	BX	46	VAL
45	BY	9	THR
45	BY	51	SER
46	BZ	4	ASP
46	BZ	7	PRO
46	BZ	8	LYS
46	BZ	26	SER
46	BZ	60	PHE
47	B0	29	VAL
47	B0	35	GLU
48	B1	45	HIS
48	B1	46	VAL
49	B2	4	THR
49	B2	43	THR
49	B2	44	VAL
50	B3	16	THR
50	B3	23	HIS
50	B3	57	VAL
50	B3	60	CYS
51	B4	5	ALA

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Mol	Chain	Res	Type
51	B4	17	VAL
51	B4	24	ARG
51	B4	29	ALA
51	B4	35	GLN
52	BI	18	ASN
2	CC	54	ILE
3	CD	24	VAL
3	CD	49	ASP
3	CD	121	ALA
4	CE	23	THR
5	CF	98	GLU
6	CG	3	ARG
6	CG	62	GLU
7	CH	66	GLN
7	CH	82	LEU
9	CJ	34	ALA
9	CJ	57	VAL
9	CJ	61	ALA
9	CJ	67	ILE
11	CL	10	PRO
11	CL	32	VAL
11	CL	33	CYS
11	CL	43	LYS
11	CL	88	ASP
11	CL	101	LEU
11	CL	111	GLN
12	CM	22	TYR
12	CM	105	ALA
13	CN	29	ILE
13	CN	50	LEU
14	CO	73	ASP
15	CP	28	ARG
18	CS	29	PRO
18	CS	62	THR
19	CT	3	ILE
19	CT	4	LYS
20	CB	22	TRP
20	CB	73	ARG
20	CB	188	THR
20	CB	206	ILE
20	CB	208	ALA
25	DC	28	PRO

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Mol	Chain	Res	Type
25	DC	29	PHE
25	DC	31	PRO
25	DC	32	LEU
25	DC	47	ARG
25	DC	63	ILE
25	DC	87	SER
25	DC	88	ALA
25	DC	97	ASP
25	DC	121	ALA
25	DC	125	PRO
25	DC	135	PRO
25	DC	149	LYS
25	DC	157	ALA
25	DC	163	ILE
25	DC	164	VAL
25	DC	176	ARG
25	DC	194	VAL
25	DC	200	MET
25	DC	246	PRO
25	DC	259	ASN
25	DC	266	ILE
26	DD	9	VAL
26	DD	34	VAL
26	DD	39	ASP
26	DD	45	TYR
26	DD	55	LYS
26	DD	82	PHE
26	DD	84	LEU
26	DD	85	ALA
26	DD	129	THR
26	DD	140	HIS
26	DD	152	PRO
26	DD	155	VAL
26	DD	157	LYS
26	DD	160	LYS
27	DE	17	THR
27	DE	45	ALA
27	DE	86	ALA
27	DE	91	ASP
27	DE	147	LEU
27	DE	162	ARG
28	DF	46	LYS

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Mol	Chain	Res	Type
28	DF	73	VAL
28	DF	84	ILE
28	DF	133	GLU
28	DF	173	ASP
29	DG	2	ARG
29	DG	4	ALA
29	DG	6	ALA
29	DG	101	VAL
30	DH	11	ASN
30	DH	38	PRO
30	DH	121	VAL
31	DJ	4	PHE
31	DJ	11	VAL
31	DJ	43	GLU
31	DJ	44	TYR
31	DJ	64	VAL
31	DJ	78	THR
31	DJ	81	ILE
31	DJ	124	VAL
32	DK	71	ARG
32	DK	72	PRO
32	DK	113	MET
32	DK	120	PRO
33	DL	7	SER
33	DL	8	PRO
33	DL	17	LYS
33	DL	38	GLN
33	DL	51	GLU
33	DL	54	GLN
33	DL	72	ALA
33	DL	92	LEU
33	DL	111	ILE
33	DL	119	PRO
33	DL	126	ARG
33	DL	127	VAL
33	DL	142	ILE
34	DM	5	LYS
34	DM	16	ARG
34	DM	17	ASN
34	DM	65	ILE
34	DM	71	LYS
34	DM	72	PRO

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Mol	Chain	Res	Type
34	DM	84	LYS
34	DM	86	LYS
34	DM	101	VAL
34	DM	106	ASP
34	DM	109	PRO
35	DN	91	ALA
36	DO	23	ALA
36	DO	27	VAL
36	DO	54	VAL
36	DO	63	LYS
36	DO	90	VAL
37	DP	17	PRO
37	DP	23	ASP
37	DP	24	THR
37	DP	41	ALA
37	DP	60	VAL
37	DP	71	ARG
37	DP	72	VAL
37	DP	76	HIS
37	DP	77	SER
37	DP	79	VAL
37	DP	80	VAL
37	DP	82	SER
37	DP	105	LYS
37	DP	110	LYS
37	DP	111	GLU
38	DQ	73	ILE
38	DQ	89	ILE
38	DQ	93	ILE
39	DR	22	LEU
39	DR	27	ILE
39	DR	73	LYS
39	DR	75	VAL
39	DR	89	HIS
39	DR	90	ARG
39	DR	97	LYS
40	DS	60	HIS
40	DS	103	ILE
41	DT	27	SER
41	DT	57	VAL
41	DT	62	VAL
41	DT	95	PHE

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Mol	Chain	Res	Type
42	DU	30	SER
42	DU	63	ALA
42	DU	66	VAL
42	DU	75	ALA
43	DW	30	VAL
43	DW	34	SER
43	DW	44	PHE
43	DW	57	THR
43	DW	58	LEU
43	DW	66	VAL
46	DZ	21	VAL
46	DZ	28	VAL
46	DZ	31	ASP
46	DZ	40	CYS
47	D0	31	LYS
47	D0	45	ASP
48	D1	14	ALA
48	D1	23	THR
48	D1	46	VAL
49	D2	44	VAL
50	D3	30	HIS
51	D4	3	VAL
51	D4	17	VAL
51	D4	19	ARG
51	D4	24	ARG
51	D4	26	ILE
51	D4	27	CYS
52	DI	18	ASN
2	AC	100	ILE
2	AC	153	SER
3	AD	31	CYS
3	AD	107	GLY
3	AD	152	SER
3	AD	172	VAL
3	AD	192	ALA
4	AE	20	VAL
4	AE	76	ASN
4	AE	107	GLY
5	AF	62	MET
6	AG	18	GLY
6	AG	129	ASN
8	AI	43	ALA

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Mol	Chain	Res	Type
8	AI	57	VAL
8	AI	108	ARG
9	AJ	74	VAL
9	AJ	75	ASP
10	AK	53	GLY
11	AL	10	PRO
11	AL	84	GLY
12	AM	3	ILE
12	AM	7	ASN
12	AM	66	GLY
12	AM	104	ASN
13	AN	51	PRO
13	AN	52	ARG
13	AN	61	ASN
13	AN	94	GLY
15	AP	29	ASN
16	AQ	6	THR
16	AQ	34	GLY
16	AQ	69	THR
17	AR	20	ILE
17	AR	22	TYR
17	AR	71	ASP
18	AS	5	LYS
19	AT	3	ILE
19	AT	85	LEU
21	AU	32	ARG
21	AU	33	ARG
21	AU	52	VAL
24	BV	45	ASP
25	BC	4	LYS
25	BC	14	HIS
25	BC	48	ILE
25	BC	51	ARG
25	BC	90	ILE
25	BC	107	LYS
25	BC	111	ALA
25	BC	143	VAL
25	BC	157	ALA
25	BC	163	ILE
25	BC	194	VAL
25	BC	208	GLY
25	BC	209	ALA

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Mol	Chain	Res	Type
25	BC	226	PRO
26	BD	2	ILE
26	BD	16	THR
26	BD	27	ILE
26	BD	30	GLU
26	BD	42	ASN
26	BD	101	PHE
26	BD	104	VAL
26	BD	125	TRP
26	BD	129	THR
26	BD	169	ARG
26	BD	202	ILE
26	BD	205	PRO
26	BD	206	ALA
26	BD	207	VAL
27	BE	23	PHE
27	BE	26	ALA
27	BE	37	ALA
27	BE	47	LYS
27	BE	49	ARG
27	BE	54	GLY
27	BE	61	ARG
27	BE	68	ALA
27	BE	86	ALA
27	BE	90	GLN
27	BE	104	ALA
27	BE	120	VAL
27	BE	128	ALA
27	BE	146	VAL
27	BE	148	ILE
27	BE	169	VAL
27	BE	171	ASP
28	BF	14	LYS
28	BF	114	ARG
28	BF	136	ILE
28	BF	153	ILE
28	BF	162	ASP
28	BF	176	PHE
29	BG	2	ARG
29	BG	32	LEU
29	BG	42	VAL
29	BG	47	ASN

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Mol	Chain	Res	Type
29	BG	173	ALA
30	BH	8	LYS
30	BH	44	ILE
30	BH	65	ALA
30	BH	127	GLU
30	BH	134	VAL
31	BJ	6	ALA
31	BJ	18	VAL
31	BJ	79	GLY
31	BJ	81	ILE
31	BJ	84	ILE
31	BJ	138	GLN
32	BK	17	ARG
32	BK	92	GLU
33	BL	15	ALA
33	BL	29	LYS
33	BL	31	GLY
33	BL	53	GLY
33	BL	94	THR
33	BL	95	LEU
33	BL	117	THR
33	BL	120	VAL
33	BL	143	GLU
34	BM	7	THR
34	BM	36	VAL
34	BM	58	LYS
34	BM	70	ASP
34	BM	86	LYS
34	BM	87	GLY
34	BM	104	GLU
34	BM	132	THR
35	BN	11	ASN
35	BN	94	TYR
35	BN	104	ALA
35	BN	114	GLU
36	BO	9	ARG
36	BO	21	LEU
36	BO	28	VAL
36	BO	38	GLN
36	BO	47	VAL
36	BO	49	VAL
36	BO	95	SER

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Mol	Chain	Res	Type
36	BO	115	LEU
37	BP	29	VAL
37	BP	30	TRP
37	BP	41	ALA
37	BP	56	SER
37	BP	59	THR
37	BP	62	LYS
37	BP	72	VAL
37	BP	76	HIS
37	BP	80	VAL
37	BP	81	ASP
37	BP	97	TYR
37	BP	102	ARG
37	BP	105	LYS
37	BP	110	LYS
37	BP	112	ARG
38	BQ	21	LYS
38	BQ	23	TYR
38	BQ	33	VAL
38	BQ	78	PHE
38	BQ	91	ARG
38	BQ	93	ILE
39	BR	17	GLY
39	BR	25	LEU
39	BR	29	THR
39	BR	37	GLU
39	BR	87	GLN
39	BR	98	ILE
39	BR	100	GLY
40	BS	3	THR
40	BS	5	ALA
40	BS	29	VAL
40	BS	67	ASP
40	BS	96	ILE
41	BT	4	GLU
41	BT	6	ARG
41	BT	14	PRO
41	BT	15	HIS
41	BT	30	ILE
41	BT	34	VAL
41	BT	67	VAL
41	BT	74	ILE

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Mol	Chain	Res	Type
41	BT	81	LYS
41	BT	85	VAL
42	BU	14	THR
42	BU	24	VAL
42	BU	35	VAL
42	BU	45	GLN
42	BU	54	PRO
42	BU	60	LYS
42	BU	64	ILE
42	BU	71	ILE
42	BU	77	GLY
43	BW	23	LYS
43	BW	38	ARG
43	BW	40	ARG
43	BW	50	VAL
43	BW	63	ASP
43	BW	72	GLY
43	BW	77	LYS
44	BX	9	LYS
44	BX	16	THR
44	BX	36	GLN
44	BX	40	SER
45	BY	3	THR
45	BY	12	ALA
46	BZ	9	TYR
46	BZ	39	LYS
46	BZ	40	CYS
46	BZ	58	ASP
47	B0	41	HIS
47	B0	49	ARG
49	B2	36	ALA
50	B3	15	LYS
51	B4	20	ASP
51	B4	22	VAL
2	CC	25	THR
2	CC	63	ILE
2	CC	65	VAL
2	CC	95	GLY
2	CC	128	MET
2	CC	153	SER
3	CD	14	GLU
3	CD	47	LEU

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Mol	Chain	Res	Type
3	CD	120	LYS
3	CD	169	TRP
4	CE	15	ILE
4	CE	17	VAL
4	CE	107	GLY
5	CF	89	VAL
6	CG	16	LYS
6	CG	113	LYS
6	CG	130	LYS
7	CH	69	ALA
7	CH	122	GLY
8	CI	24	ASN
8	CI	57	VAL
9	CJ	33	GLY
9	CJ	37	ARG
9	CJ	60	ASP
9	CJ	74	VAL
9	CJ	77	VAL
9	CJ	94	ALA
10	CK	14	GLN
10	CK	17	ASP
11	CL	13	ARG
11	CL	97	VAL
11	CL	104	SER
11	CL	117	GLY
11	CL	121	PRO
12	CM	3	ILE
12	CM	7	ASN
12	CM	23	GLY
13	CN	21	ALA
13	CN	51	PRO
14	CO	3	SER
15	CP	46	LYS
15	CP	79	ASN
16	CQ	28	VAL
16	CQ	47	ASP
16	CQ	56	ASP
16	CQ	69	THR
17	CR	21	ASP
18	CS	27	LYS
18	CS	69	LYS
20	CB	18	GLN

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Mol	Chain	Res	Type
20	CB	19	THR
20	CB	72	LYS
20	CB	97	GLY
20	CB	124	THR
20	CB	193	ASP
21	CU	22	CYS
24	DV	14	LYS
24	DV	45	ASP
24	DV	70	ILE
25	DC	21	PRO
25	DC	22	GLU
25	DC	39	SER
25	DC	43	ASN
25	DC	52	HIS
25	DC	53	ILE
25	DC	65	ASP
25	DC	70	LYS
25	DC	90	ILE
25	DC	93	VAL
25	DC	106	PRO
25	DC	156	SER
25	DC	173	LEU
25	DC	199	HIS
25	DC	201	LEU
25	DC	210	ALA
25	DC	227	VAL
25	DC	252	LYS
26	DD	43	ASP
26	DD	73	VAL
26	DD	77	ARG
26	DD	89	GLU
26	DD	95	SER
26	DD	145	SER
26	DD	168	GLU
26	DD	184	ARG
26	DD	185	ASN
26	DD	187	LEU
26	DD	193	VAL
26	DD	197	THR
26	DD	205	PRO
26	DD	206	ALA
27	DE	4	VAL

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Mol	Chain	Res	Type
27	DE	7	ASP
27	DE	13	THR
27	DE	43	THR
27	DE	48	THR
27	DE	49	ARG
27	DE	54	GLY
27	DE	59	PRO
27	DE	84	THR
27	DE	85	PHE
27	DE	110	SER
27	DE	115	GLN
27	DE	126	VAL
27	DE	148	ILE
27	DE	149	ILE
27	DE	153	LEU
27	DE	166	LYS
27	DE	168	ASP
27	DE	195	GLN
28	DF	20	ASN
28	DF	49	LEU
28	DF	66	ILE
28	DF	80	GLN
28	DF	83	PRO
28	DF	109	ARG
28	DF	136	ILE
28	DF	149	ARG
28	DF	150	GLY
28	DF	174	PHE
29	DG	5	LYS
29	DG	37	ASN
30	DH	10	ALA
30	DH	25	TYR
30	DH	30	LEU
30	DH	82	SER
30	DH	91	PHE
31	DJ	50	THR
31	DJ	65	THR
31	DJ	128	ASN
31	DJ	129	GLU
31	DJ	131	ASN
31	DJ	135	GLN
32	DK	3	GLN

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Mol	Chain	Res	Type
32	DK	10	VAL
32	DK	18	ARG
32	DK	83	ALA
33	DL	15	ALA
33	DL	20	GLY
33	DL	27	LEU
33	DL	40	SER
33	DL	64	PHE
33	DL	108	ALA
33	DL	109	LYS
33	DL	110	VAL
34	DM	80	VAL
34	DM	102	LEU
35	DN	100	CYS
35	DN	116	VAL
36	DO	22	GLY
36	DO	65	THR
36	DO	84	GLU
36	DO	96	GLY
36	DO	99	TYR
36	DO	101	GLY
37	DP	27	VAL
37	DP	49	ILE
37	DP	69	VAL
37	DP	75	THR
37	DP	83	ILE
37	DP	91	VAL
38	DQ	5	ARG
38	DQ	30	VAL
38	DQ	31	TYR
38	DQ	85	ALA
38	DQ	88	GLU
39	DR	4	VAL
39	DR	5	PHE
39	DR	11	GLN
39	DR	59	ILE
39	DR	81	LYS
39	DR	82	HIS
39	DR	93	PHE
40	DS	4	ILE
40	DS	12	SER
40	DS	26	GLY

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Mol	Chain	Res	Type
40	DS	41	LYS
40	DS	64	ALA
40	DS	69	LEU
40	DS	87	PRO
41	DT	7	LEU
41	DT	15	HIS
41	DT	31	VAL
41	DT	36	LYS
41	DT	52	GLU
41	DT	63	VAL
41	DT	71	GLY
41	DT	85	VAL
42	DU	4	ILE
42	DU	9	GLU
42	DU	24	VAL
42	DU	28	LEU
42	DU	42	LYS
42	DU	93	ARG
43	DW	4	LYS
43	DW	10	ARG
43	DW	18	LYS
43	DW	45	HIS
43	DW	62	ALA
43	DW	75	ASN
43	DW	77	LYS
44	DX	20	ASN
45	DY	12	ALA
46	DZ	3	LYS
46	DZ	10	GLU
47	D0	35	GLU
47	D0	47	TYR
48	D1	30	PRO
48	D1	38	PHE
48	D1	47	ILE
48	D1	50	GLU
49	D2	5	PHE
49	D2	18	PHE
49	D2	25	LYS
50	D3	5	THR
50	D3	26	ALA
51	D4	10	LEU
2	AC	25	THR

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Mol	Chain	Res	Type
2	AC	83	VAL
2	AC	136	ALA
2	AC	180	ASP
3	AD	28	ASP
3	AD	148	ALA
3	AD	151	GLN
4	AE	43	GLY
5	AF	65	GLU
6	AG	75	LYS
7	AH	46	GLU
7	AH	65	PHE
9	AJ	61	ALA
9	AJ	62	ARG
9	AJ	92	LEU
9	AJ	93	ALA
10	AK	77	GLY
10	AK	125	LYS
11	AL	13	ARG
11	AL	15	VAL
11	AL	47	ALA
11	AL	120	ARG
12	AM	22	TYR
12	AM	98	GLY
13	AN	2	LYS
13	AN	80	ARG
17	AR	32	ILE
18	AS	3	SER
20	AB	18	GLN
20	AB	20	ARG
20	AB	27	LYS
20	AB	76	SER
20	AB	88	GLN
20	AB	186	VAL
21	AU	24	LYS
21	AU	35	GLU
21	AU	40	PRO
24	BV	54	ALA
24	BV	64	VAL
24	BV	71	LYS
25	BC	34	GLU
25	BC	41	GLY
25	BC	52	HIS

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Mol	Chain	Res	Type
25	BC	53	ILE
25	BC	58	LYS
25	BC	59	GLN
25	BC	65	ASP
25	BC	66	PHE
25	BC	91	ALA
25	BC	93	VAL
25	BC	108	GLY
25	BC	125	PRO
25	BC	231	HIS
25	BC	239	PHE
25	BC	251	THR
25	BC	255	LYS
25	BC	259	ASN
26	BD	18	ASP
26	BD	21	SER
26	BD	23	PRO
26	BD	105	LYS
26	BD	189	VAL
27	BE	11	ALA
27	BE	24	ASN
27	BE	62	GLN
27	BE	93	SER
27	BE	191	ASP
28	BF	19	PHE
28	BF	34	THR
28	BF	67	THR
28	BF	75	GLY
28	BF	81	GLY
28	BF	88	VAL
28	BF	173	ASP
29	BG	127	GLN
30	BH	125	THR
31	BJ	50	THR
31	BJ	74	TYR
31	BJ	83	GLY
31	BJ	85	LYS
32	BK	46	ALA
32	BK	112	PHE
33	BL	22	GLY
33	BL	27	LEU
33	BL	38	GLN

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Mol	Chain	Res	Type
33	BL	48	ARG
33	BL	57	LEU
34	BM	35	ALA
34	BM	59	ARG
34	BM	77	PRO
34	BM	105	MET
34	BM	127	LYS
35	BN	6	SER
35	BN	102	PHE
36	BO	32	PRO
36	BO	52	SER
36	BO	68	LYS
37	BP	17	PRO
37	BP	48	ALA
37	BP	52	ARG
37	BP	71	ARG
37	BP	78	PRO
38	BQ	73	ILE
39	BR	18	GLN
39	BR	42	ALA
39	BR	45	GLU
39	BR	73	LYS
39	BR	76	LYS
39	BR	77	PHE
40	BS	13	SER
40	BS	69	LEU
41	BT	13	ALA
41	BT	61	LEU
42	BU	18	LYS
42	BU	39	ASN
42	BU	67	SER
42	BU	72	PHE
42	BU	96	LYS
43	BW	6	GLY
43	BW	49	ASN
43	BW	62	ALA
43	BW	65	LYS
43	BW	68	PHE
44	BX	4	LYS
44	BX	28	LEU
45	BY	14	GLY
45	BY	16	LEU

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Mol	Chain	Res	Type
45	BY	42	ALA
47	B0	4	GLN
48	B1	8	ILE
48	B1	35	LEU
48	B1	43	ARG
51	B4	3	VAL
51	B4	10	LEU
52	BI	14	ALA
52	BI	23	VAL
2	CC	3	LYS
2	CC	145	ALA
2	CC	167	TYR
5	CF	85	ILE
5	CF	95	ALA
6	CG	55	LYS
7	CH	30	LYS
7	CH	53	ASP
8	CI	34	LEU
8	CI	55	ASP
8	CI	106	ASP
9	CJ	68	ARG
9	CJ	93	ALA
9	CJ	100	ILE
10	CK	52	ARG
10	CK	53	GLY
10	CK	88	PRO
10	CK	108	ASN
11	CL	24	GLU
11	CL	73	LEU
12	CM	62	PHE
12	CM	65	GLU
13	CN	33	VAL
13	CN	42	ASN
13	CN	43	ALA
13	CN	52	ARG
13	CN	69	PRO
13	CN	70	HIS
15	CP	47	GLU
18	CS	8	PRO
18	CS	34	SER
18	CS	63	ASP
19	CT	68	LYS

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Mol	Chain	Res	Type
20	CB	43	GLU
20	CB	44	LYS
20	CB	94	ARG
20	CB	95	TRP
20	CB	114	LYS
20	CB	128	LEU
20	CB	204	ASP
21	CU	23	GLU
25	DC	66	PHE
25	DC	115	ILE
25	DC	172	THR
25	DC	184	GLU
25	DC	243	PRO
25	DC	265	PHE
26	DD	47	ALA
26	DD	139	SER
26	DD	141	ARG
26	DD	147	GLY
26	DD	174	SER
27	DE	30	GLN
27	DE	38	GLY
27	DE	67	ARG
27	DE	90	GLN
27	DE	139	LYS
27	DE	172	ALA
27	DE	190	ALA
28	DF	42	ALA
28	DF	43	ILE
28	DF	88	VAL
28	DF	104	THR
28	DF	134	GLN
28	DF	138	PRO
28	DF	176	PHE
29	DG	11	PRO
29	DG	45	ALA
29	DG	60	GLY
29	DG	173	ALA
30	DH	31	VAL
30	DH	37	VAL
30	DH	86	ASP
30	DH	92	GLY
31	DJ	8	PRO

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Mol	Chain	Res	Type
31	DJ	41	LYS
31	DJ	53	TYR
31	DJ	58	ASN
31	DJ	75	TYR
31	DJ	77	HIS
31	DJ	85	LYS
31	DJ	137	PRO
32	DK	12	ASP
32	DK	17	ARG
33	DL	14	LYS
33	DL	44	GLY
33	DL	45	GLY
33	DL	52	GLY
33	DL	55	MET
33	DL	116	VAL
33	DL	141	LYS
34	DM	4	PRO
34	DM	127	LYS
34	DM	128	THR
35	DN	32	GLU
36	DO	34	HIS
36	DO	38	GLN
36	DO	87	ILE
37	DP	39	LEU
37	DP	97	TYR
37	DP	112	ARG
38	DQ	53	LYS
38	DQ	86	SER
39	DR	3	ALA
39	DR	10	LYS
39	DR	42	ALA
39	DR	64	VAL
39	DR	86	GLN
40	DS	9	HIS
40	DS	99	ARG
41	DT	16	VAL
41	DT	39	THR
41	DT	68	LYS
41	DT	96	VAL
42	DU	62	ALA
42	DU	64	ILE
42	DU	99	SER

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Mol	Chain	Res	Type
43	DW	5	ALA
43	DW	9	THR
43	DW	65	LYS
44	DX	28	LEU
44	DX	31	GLN
44	DX	35	GLY
44	DX	52	ARG
45	DY	37	ARG
46	DZ	26	SER
47	D0	36	LYS
47	D0	43	THR
48	D1	32	LYS
49	D2	7	PRO
49	D2	8	SER
50	D3	10	ALA
50	D3	47	ALA
51	D4	11	CYS
51	D4	31	PRO
52	DI	23	VAL
2	AC	19	SER
2	AC	54	ILE
2	AC	65	VAL
3	AD	7	LYS
3	AD	27	ILE
3	AD	29	THR
3	AD	143	SER
4	AE	34	ALA
4	AE	128	GLY
5	AF	64	VAL
5	AF	94	HIS
6	AG	84	TYR
8	AI	44	ARG
8	AI	55	ASP
8	AI	67	LYS
8	AI	127	SER
9	AJ	36	VAL
9	AJ	56	HIS
9	AJ	95	GLY
11	AL	14	LYS
11	AL	122	LYS
13	AN	21	ALA
15	AP	27	ALA

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Mol	Chain	Res	Type
15	AP	42	ILE
15	AP	54	LEU
18	AS	8	PRO
18	AS	15	LEU
18	AS	39	ILE
19	AT	59	ARG
20	AB	41	ASN
20	AB	150	ILE
20	AB	200	PRO
20	AB	205	ALA
21	AU	6	ARG
21	AU	36	PHE
21	AU	37	TYR
25	BC	50	THR
25	BC	61	TYR
25	BC	151	GLY
25	BC	165	ALA
25	BC	236	GLY
25	BC	257	ARG
25	BC	258	SER
25	BC	265	PHE
26	BD	47	ALA
26	BD	87	GLY
26	BD	90	PHE
26	BD	145	SER
26	BD	185	ASN
26	BD	190	LYS
26	BD	193	VAL
27	BE	39	ALA
27	BE	53	THR
27	BE	106	LYS
28	BF	7	TYR
28	BF	45	ASP
28	BF	74	ALA
28	BF	172	PHE
29	BG	48	THR
29	BG	155	PRO
29	BG	174	LYS
30	BH	20	ASN
31	BJ	25	LEU
31	BJ	36	LEU
31	BJ	97	PRO

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Mol	Chain	Res	Type
31	BJ	127	GLY
32	BK	6	THR
32	BK	121	GLU
33	BL	42	SER
33	BL	54	GLN
33	BL	82	LEU
33	BL	109	LYS
34	BM	10	ARG
34	BM	28	PHE
34	BM	133	LYS
36	BO	36	TYR
36	BO	87	ILE
36	BO	90	VAL
37	BP	77	SER
39	BR	27	ILE
39	BR	95	ASP
40	BS	11	ARG
40	BS	99	ARG
41	BT	5	GLU
42	BU	3	LYS
42	BU	59	GLU
43	BW	32	ALA
44	BX	7	ARG
45	BY	15	ARG
45	BY	17	PRO
46	BZ	23	LYS
47	B0	5	ASN
48	B1	5	ARG
48	B1	16	THR
48	B1	38	PHE
48	B1	47	ILE
49	B2	42	LEU
51	B4	7	VAL
52	BI	49	GLU
52	BI	64	ARG
2	CC	50	SER
2	CC	59	PRO
2	CC	60	ALA
2	CC	110	LEU
2	CC	144	GLY
3	CD	22	SER
3	CD	26	ALA

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Mol	Chain	Res	Type
3	CD	119	HIS
3	CD	166	LYS
4	CE	87	VAL
4	CE	146	MET
6	CG	66	GLU
6	CG	78	ARG
6	CG	128	GLU
6	CG	129	ASN
7	CH	94	VAL
8	CI	58	GLU
8	CI	90	ASP
9	CJ	56	HIS
9	CJ	75	ASP
11	CL	23	LEU
12	CM	97	ARG
12	CM	102	LYS
12	CM	104	ASN
13	CN	62	ARG
13	CN	65	GLN
15	CP	27	ALA
18	CS	5	LYS
18	CS	22	VAL
18	CS	61	VAL
18	CS	77	ARG
19	CT	67	HIS
20	CB	15	PHE
20	CB	86	CYS
20	CB	126	ASP
21	CU	12	ASP
24	DV	16	ALA
24	DV	71	LYS
25	DC	12	ARG
25	DC	103	ILE
25	DC	113	ASP
25	DC	124	LYS
25	DC	152	GLN
25	DC	226	PRO
25	DC	239	PHE
26	DD	87	GLY
26	DD	173	GLN
27	DE	50	ALA
27	DE	72	SER

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Mol	Chain	Res	Type
27	DE	111	GLU
28	DF	34	THR
28	DF	36	ASN
28	DF	67	THR
28	DF	70	ARG
29	DG	36	LEU
30	DH	114	GLU
31	DJ	23	LYS
31	DJ	60	ASP
31	DJ	134	ALA
32	DK	29	HIS
32	DK	46	ALA
32	DK	84	CYS
32	DK	86	LEU
32	DK	89	ASN
33	DL	5	THR
33	DL	60	ARG
33	DL	82	LEU
33	DL	118	THR
34	DM	107	GLY
35	DN	99	LYS
35	DN	102	PHE
35	DN	113	ILE
36	DO	2	ASP
37	DP	15	ASP
37	DP	29	VAL
37	DP	33	GLU
37	DP	62	LYS
37	DP	73	PHE
37	DP	90	ALA
38	DQ	4	LYS
38	DQ	9	ALA
39	DR	71	LYS
39	DR	87	GLN
39	DR	98	ILE
40	DS	8	ARG
40	DS	13	SER
40	DS	23	LEU
40	DS	27	LYS
40	DS	80	PRO
41	DT	32	LEU
41	DT	67	VAL

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Mol	Chain	Res	Type
41	DT	79	ASP
42	DU	25	LYS
42	DU	58	VAL
42	DU	94	PHE
43	DW	16	GLU
43	DW	22	VAL
43	DW	61	LYS
43	DW	63	ASP
43	DW	76	ARG
44	DX	36	GLN
45	DY	51	SER
46	DZ	32	LEU
46	DZ	35	ASP
47	D0	20	ALA
47	D0	51	ARG
48	D1	5	ARG
48	D1	43	ARG
48	D1	51	ALA
49	D2	17	GLY
49	D2	22	MET
51	D4	9	LYS
52	DI	14	ALA
2	AC	14	VAL
2	AC	59	PRO
2	AC	145	ALA
2	AC	186	SER
3	AD	6	PRO
3	AD	43	ARG
3	AD	177	MET
6	AG	57	GLU
6	AG	137	ARG
10	AK	106	ILE
10	AK	127	ARG
13	AN	14	ALA
17	AR	47	ARG
18	AS	4	LEU
18	AS	27	LYS
18	AS	66	VAL
20	AB	87	ASP
20	AB	121	GLN
21	AU	13	VAL
21	AU	26	GLY

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Mol	Chain	Res	Type
24	BV	70	ILE
24	BV	93	ARG
25	BC	30	ALA
25	BC	40	GLY
25	BC	112	GLY
25	BC	115	ILE
25	BC	196	ASN
25	BC	217	PRO
25	BC	230	PRO
26	BD	52	THR
26	BD	162	ALA
26	BD	174	SER
28	BF	10	GLU
28	BF	22	ASN
28	BF	77	LYS
28	BF	87	LYS
28	BF	122	ASP
29	BG	11	PRO
29	BG	43	LYS
29	BG	159	LYS
31	BJ	46	PRO
31	BJ	53	TYR
31	BJ	66	GLY
32	BK	3	GLN
32	BK	4	GLU
32	BK	85	VAL
33	BL	114	GLY
33	BL	125	LEU
33	BL	128	THR
34	BM	29	GLY
34	BM	30	SER
34	BM	125	PRO
35	BN	73	ASN
36	BO	31	THR
37	BP	39	LEU
38	BQ	20	ALA
38	BQ	26	ALA
39	BR	72	VAL
40	BS	9	HIS
40	BS	57	ASN
40	BS	64	ALA
40	BS	71	VAL

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Mol	Chain	Res	Type
41	BT	39	THR
41	BT	88	LYS
43	BW	5	ALA
43	BW	31	LEU
45	BY	10	ARG
46	BZ	5	ILE
47	B0	53	VAL
48	B1	21	THR
49	B2	6	GLN
49	B2	17	GLY
50	B3	24	LYS
50	B3	31	ILE
2	CC	100	ILE
2	CC	129	PHE
3	CD	191	SER
3	CD	203	TYR
3	CD	204	SER
5	CF	53	LYS
5	CF	68	GLN
8	CI	8	THR
8	CI	9	GLY
8	CI	121	ARG
9	CJ	36	VAL
9	CJ	86	ALA
10	CK	49	SER
12	CM	21	ILE
13	CN	31	SER
13	CN	73	LEU
13	CN	74	ARG
15	CP	44	SER
15	CP	55	ASP
15	CP	69	ASP
17	CR	63	TYR
20	CB	211	LEU
21	CU	39	LYS
25	DC	50	THR
25	DC	58	LYS
25	DC	137	GLY
25	DC	175	LEU
26	DD	2	ILE
26	DD	49	GLN
26	DD	149	ASN

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Mol	Chain	Res	Type
27	DE	18	THR
27	DE	46	GLN
27	DE	89	PRO
27	DE	157	LEU
27	DE	188	MET
27	DE	192	ALA
28	DF	10	GLU
28	DF	153	ILE
29	DG	44	HIS
29	DG	151	ARG
29	DG	164	ALA
30	DH	66	ASN
31	DJ	111	LYS
31	DJ	125	TYR
31	DJ	139	VAL
32	DK	14	SER
32	DK	26	GLY
32	DK	85	VAL
33	DL	74	THR
34	DM	77	PRO
34	DM	82	MET
35	DN	41	ALA
35	DN	107	ASN
36	DO	18	LEU
36	DO	103	VAL
37	DP	59	THR
37	DP	106	ALA
38	DQ	78	PHE
38	DQ	87	VAL
39	DR	37	GLU
39	DR	84	ARG
40	DS	22	ASP
40	DS	92	ARG
41	DT	60	THR
41	DT	66	LYS
41	DT	86	THR
42	DU	57	ILE
43	DW	55	ASP
44	DX	51	ALA
44	DX	61	ALA
46	DZ	19	GLY
48	D1	21	THR

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Mol	Chain	Res	Type
48	D1	29	LYS
49	D2	14	ARG
49	D2	15	SER
49	D2	45	SER
50	D3	13	PHE
50	D3	29	ARG
50	D3	33	THR
51	D4	8	LYS
52	DI	5	GLN
4	AE	25	LYS
6	AG	6	ILE
6	AG	86	VAL
7	AH	82	LEU
10	AK	91	GLY
12	AM	65	GLU
13	AN	33	VAL
16	AQ	28	VAL
18	AS	36	ARG
18	AS	52	ASN
18	AS	67	GLY
21	AU	23	GLU
25	BC	164	VAL
25	BC	224	MET
26	BD	19	GLY
26	BD	99	GLU
26	BD	180	VAL
27	BE	75	SER
27	BE	145	ASP
27	BE	186	VAL
28	BF	83	PRO
28	BF	175	PRO
29	BG	84	LYS
30	BH	108	VAL
31	BJ	121	LYS
32	BK	49	ARG
34	BM	21	ALA
34	BM	26	VAL
34	BM	97	GLN
35	BN	115	LEU
37	BP	26	GLU
39	BR	7	SER
39	BR	12	HIS

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Mol	Chain	Res	Type
42	BU	56	GLY
43	BW	42	THR
46	BZ	32	LEU
46	BZ	66	ILE
47	B0	34	GLY
50	B3	26	ALA
2	CC	127	VAL
3	CD	35	GLN
3	CD	37	PRO
4	CE	20	VAL
4	CE	78	GLY
5	CF	84	VAL
5	CF	99	ALA
6	CG	99	ALA
8	CI	23	GLY
8	CI	71	ILE
9	CJ	32	THR
9	CJ	62	ARG
11	CL	99	GLY
11	CL	103	CYS
13	CN	3	GLN
15	CP	54	LEU
17	CR	36	GLY
19	CT	44	ALA
20	CB	20	ARG
20	CB	205	ALA
25	DC	60	ALA
25	DC	208	GLY
25	DC	217	PRO
25	DC	224	MET
25	DC	244	VAL
26	DD	29	VAL
26	DD	48	ILE
26	DD	53	GLY
27	DE	56	GLY
27	DE	66	GLY
27	DE	167	VAL
31	DJ	46	PRO
31	DJ	67	ASN
33	DL	9	ALA
33	DL	49	GLY
35	DN	123	GLU

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Mol	Chain	Res	Type
37	DP	26	GLU
39	DR	54	VAL
39	DR	63	VAL
39	DR	102	SER
40	DS	3	THR
40	DS	42	LYS
41	DT	76	ARG
41	DT	92	ASN
42	DU	65	GLN
43	DW	52	CYS
44	DX	8	GLU
50	D3	49	VAL
51	D4	28	SER
3	AD	175	GLY
11	AL	121	PRO
14	AO	26	VAL
20	AB	24	PRO
26	BD	96	ILE
27	BE	82	GLY
29	BG	7	PRO
33	BL	11	GLY
33	BL	102	GLY
40	BS	4	ILE
43	BW	35	ILE
47	B0	46	GLY
6	CG	18	GLY
21	CU	13	VAL
26	DD	76	GLY
27	DE	58	LYS
30	DH	85	GLY
32	DK	103	VAL
33	DL	32	GLY
34	DM	85	GLY
36	DO	28	VAL
38	DQ	72	GLY
39	DR	33	VAL
39	DR	47	VAL
4	AE	26	GLY
17	AR	43	ILE
25	BC	16	VAL
25	BC	214	GLY
26	BD	73	VAL

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Mol	Chain	Res	Type
26	BD	203	VAL
28	BF	148	VAL
29	BG	110	HIS
30	BH	34	GLY
30	BH	95	GLY
32	BK	15	GLY
34	BM	71	LYS
34	BM	131	VAL
38	BQ	7	VAL
44	BX	11	VAL
3	CD	27	ILE
10	CK	15	VAL
12	CM	42	VAL
16	CQ	32	ILE
17	CR	20	ILE
25	DC	48	ILE
25	DC	161	VAL
30	DH	4	ILE
33	DL	100	ILE
41	DT	58	VAL
50	D3	31	ILE
50	D3	62	PRO
52	DI	34	ILE
3	AD	63	ILE
7	AH	71	VAL
8	AI	71	ILE
9	AJ	41	PRO
24	BV	84	PRO
25	BC	234	GLY
26	BD	98	VAL
27	BE	113	VAL
27	BE	187	VAL
28	BF	108	PRO
30	BH	110	VAL
36	BO	35	ILE
37	BP	69	VAL
2	CC	76	ILE
2	CC	107	LYS
4	CE	24	VAL
20	CB	123	GLY
20	CB	200	PRO
25	DC	54	GLY

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Mol	Chain	Res	Type
25	DC	195	GLY
26	DD	19	GLY
26	DD	203	VAL
33	DL	23	ILE
51	D4	23	ILE
3	AD	37	PRO
7	AH	74	ILE
8	AI	110	VAL
11	AL	62	VAL
12	AM	15	VAL
28	BF	11	VAL
29	BG	8	VAL
29	BG	99	GLY
33	BL	16	GLY
37	BP	3	ILE
8	CI	68	GLY
10	CK	119	GLY
16	CQ	4	ILE
25	DC	16	VAL
27	DE	120	VAL
28	DF	108	PRO
33	DL	89	VAL
36	DO	31	THR
36	DO	58	ILE
38	DQ	6	GLY
42	DU	11	ILE
47	D0	53	VAL
49	D2	6	GLN
8	AI	39	GLY
27	BE	76	PRO
34	BM	89	VAL
47	B0	42	ILE
29	DG	117	PRO
30	DH	32	PRO
30	DH	107	GLY

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	170/190 (90%)	149 (88%)	21 (12%)	5	26
2	CC	170/190 (90%)	150 (88%)	20 (12%)	6	28
3	AD	172/173 (99%)	145 (84%)	27 (16%)	3	17
3	CD	172/173 (99%)	143 (83%)	29 (17%)	2	13
4	AE	113/126 (90%)	96 (85%)	17 (15%)	3	19
4	CE	113/126 (90%)	94 (83%)	19 (17%)	2	14
5	AF	87/116 (75%)	68 (78%)	19 (22%)	1	5
5	CF	87/116 (75%)	76 (87%)	11 (13%)	5	26
6	AG	123/147 (84%)	106 (86%)	17 (14%)	4	22
6	CG	125/147 (85%)	111 (89%)	14 (11%)	7	30
7	AH	104/105 (99%)	92 (88%)	12 (12%)	6	29
7	CH	104/105 (99%)	93 (89%)	11 (11%)	8	34
8	AI	105/107 (98%)	90 (86%)	15 (14%)	4	21
8	CI	105/107 (98%)	91 (87%)	14 (13%)	4	24
9	AJ	86/90 (96%)	70 (81%)	16 (19%)	2	9
9	CJ	86/90 (96%)	78 (91%)	8 (9%)	10	40
10	AK	90/99 (91%)	72 (80%)	18 (20%)	1	7
10	CK	90/99 (91%)	79 (88%)	11 (12%)	6	27
11	AL	103/104 (99%)	90 (87%)	13 (13%)	5	26
11	CL	103/104 (99%)	86 (84%)	17 (16%)	2	14
12	AM	92/96 (96%)	76 (83%)	16 (17%)	2	12
12	CM	91/96 (95%)	80 (88%)	11 (12%)	6	27
13	AN	79/84 (94%)	70 (89%)	9 (11%)	7	29
13	CN	79/84 (94%)	73 (92%)	6 (8%)	15	50
14	AO	76/77 (99%)	69 (91%)	7 (9%)	11	40
14	CO	76/77 (99%)	67 (88%)	9 (12%)	6	28
15	AP	65/65 (100%)	54 (83%)	11 (17%)	2	13
15	CP	65/65 (100%)	57 (88%)	8 (12%)	5	27
16	AQ	74/78 (95%)	64 (86%)	10 (14%)	4	23
16	CQ	75/78 (96%)	68 (91%)	7 (9%)	10	40
17	AR	48/65 (74%)	44 (92%)	4 (8%)	13	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	CR	48/65 (74%)	42 (88%)	6 (12%)	5	26
18	AS	70/79 (89%)	59 (84%)	11 (16%)	3	17
18	CS	71/79 (90%)	58 (82%)	13 (18%)	2	10
19	AT	65/66 (98%)	55 (85%)	10 (15%)	3	17
19	CT	65/66 (98%)	55 (85%)	10 (15%)	3	17
20	AB	180/199 (90%)	148 (82%)	32 (18%)	2	11
20	CB	180/199 (90%)	153 (85%)	27 (15%)	3	19
21	AU	44/61 (72%)	37 (84%)	7 (16%)	3	16
21	CU	44/61 (72%)	31 (70%)	13 (30%)	0	3
24	BV	78/78 (100%)	64 (82%)	14 (18%)	2	11
24	DV	78/78 (100%)	72 (92%)	6 (8%)	15	49
25	BC	213/218 (98%)	162 (76%)	51 (24%)	1	4
25	DC	213/218 (98%)	164 (77%)	49 (23%)	1	4
26	BD	164/164 (100%)	130 (79%)	34 (21%)	1	6
26	DD	164/164 (100%)	126 (77%)	38 (23%)	1	4
27	BE	165/165 (100%)	127 (77%)	38 (23%)	1	4
27	DE	165/165 (100%)	137 (83%)	28 (17%)	2	13
28	BF	149/150 (99%)	127 (85%)	22 (15%)	3	19
28	DF	149/150 (99%)	127 (85%)	22 (15%)	3	19
29	BG	137/138 (99%)	115 (84%)	22 (16%)	3	16
29	DG	137/138 (99%)	114 (83%)	23 (17%)	2	14
30	BH	114/114 (100%)	95 (83%)	19 (17%)	2	14
30	DH	114/114 (100%)	92 (81%)	22 (19%)	1	8
31	BJ	114/116 (98%)	93 (82%)	21 (18%)	2	10
31	DJ	114/116 (98%)	92 (81%)	22 (19%)	1	8
32	BK	102/104 (98%)	81 (79%)	21 (21%)	1	6
32	DK	102/104 (98%)	89 (87%)	13 (13%)	5	25
33	BL	103/103 (100%)	76 (74%)	27 (26%)	0	3
33	DL	103/103 (100%)	78 (76%)	25 (24%)	1	4
34	BM	109/109 (100%)	81 (74%)	28 (26%)	0	3
34	DM	109/109 (100%)	75 (69%)	34 (31%)	0	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
35	BN	103/103 (100%)	84 (82%)	19 (18%)	2	10
35	DN	103/103 (100%)	87 (84%)	16 (16%)	3	17
36	BO	87/87 (100%)	68 (78%)	19 (22%)	1	5
36	DO	87/87 (100%)	73 (84%)	14 (16%)	3	16
37	BP	99/100 (99%)	78 (79%)	21 (21%)	1	6
37	DP	99/100 (99%)	72 (73%)	27 (27%)	0	3
38	BQ	89/90 (99%)	72 (81%)	17 (19%)	2	8
38	DQ	89/90 (99%)	70 (79%)	19 (21%)	1	6
39	BR	84/84 (100%)	72 (86%)	12 (14%)	4	21
39	DR	84/84 (100%)	67 (80%)	17 (20%)	1	7
40	BS	93/93 (100%)	83 (89%)	10 (11%)	7	33
40	DS	93/93 (100%)	80 (86%)	13 (14%)	4	21
41	BT	83/84 (99%)	65 (78%)	18 (22%)	1	5
41	DT	83/84 (99%)	70 (84%)	13 (16%)	3	17
42	BU	83/85 (98%)	66 (80%)	17 (20%)	1	7
42	DU	83/85 (98%)	65 (78%)	18 (22%)	1	5
43	BW	62/63 (98%)	50 (81%)	12 (19%)	1	8
43	DW	62/63 (98%)	51 (82%)	11 (18%)	2	11
44	BX	55/55 (100%)	38 (69%)	17 (31%)	0	2
44	DX	55/55 (100%)	45 (82%)	10 (18%)	2	10
45	BY	48/49 (98%)	37 (77%)	11 (23%)	1	4
45	DY	48/49 (98%)	36 (75%)	12 (25%)	1	4
46	BZ	62/62 (100%)	46 (74%)	16 (26%)	0	3
46	DZ	62/62 (100%)	54 (87%)	8 (13%)	5	25
47	B0	47/48 (98%)	36 (77%)	11 (23%)	1	4
47	D0	47/48 (98%)	38 (81%)	9 (19%)	2	8
48	B1	48/49 (98%)	35 (73%)	13 (27%)	0	3
48	D1	48/49 (98%)	38 (79%)	10 (21%)	1	6
49	B2	38/38 (100%)	29 (76%)	9 (24%)	1	4
49	D2	38/38 (100%)	29 (76%)	9 (24%)	1	4
50	B3	51/52 (98%)	41 (80%)	10 (20%)	1	8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
50	D3	51/52 (98%)	39 (76%)	12 (24%)	1	4
51	B4	34/34 (100%)	25 (74%)	9 (26%)	0	3
51	D4	34/34 (100%)	19 (56%)	15 (44%)	0	0
52	BI	109/110 (99%)	107 (98%)	2 (2%)	64	86
52	DI	109/110 (99%)	105 (96%)	4 (4%)	39	73
All	All	9341/9744 (96%)	7726 (83%)	1615 (17%)	2	12

All (1615) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	AC	31	ASN
2	AC	35	ASP
2	AC	40	GLN
2	AC	64	ARG
2	AC	82	ASP
2	AC	100	ILE
2	AC	102	ILE
2	AC	106	ARG
2	AC	113	LYS
2	AC	141	MET
2	AC	143	LEU
2	AC	146	LYS
2	AC	156	LEU
2	AC	166	TRP
2	AC	168	ARG
2	AC	174	LEU
2	AC	180	ASP
2	AC	184	ASN
2	AC	189	HIS
2	AC	190	THR
2	AC	192	TYR
3	AD	7	LYS
3	AD	8	LEU
3	AD	10	LEU
3	AD	21	LYS
3	AD	25	ARG
3	AD	35	GLN
3	AD	39	GLN
3	AD	43	ARG
3	AD	56	GLU

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Mol	Chain	Res	Type
3	AD	58	GLN
3	AD	62	ARG
3	AD	64	TYR
3	AD	69	ARG
3	AD	85	THR
3	AD	94	GLU
3	AD	147	LYS
3	AD	150	LYS
3	AD	153	ARG
3	AD	155	LYS
3	AD	158	LEU
3	AD	159	GLU
3	AD	160	LEU
3	AD	176	LYS
3	AD	177	MET
3	AD	183	ARG
3	AD	186	GLU
3	AD	194	ILE
4	AE	11	GLN
4	AE	23	THR
4	AE	25	LYS
4	AE	44	ARG
4	AE	68	ARG
4	AE	72	ASN
4	AE	81	GLN
4	AE	92	ARG
4	AE	102	THR
4	AE	110	MET
4	AE	122	VAL
4	AE	123	LEU
4	AE	125	LYS
4	AE	139	THR
4	AE	143	LEU
4	AE	151	MET
4	AE	158	LYS
5	AF	24	ARG
5	AF	37	HIS
5	AF	38	ARG
5	AF	39	LEU
5	AF	46	GLN
5	AF	53	LYS
5	AF	54	LEU

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Mol	Chain	Res	Type
5	AF	55	HIS
5	AF	61	LEU
5	AF	65	GLU
5	AF	68	GLN
5	AF	69	GLU
5	AF	71	ILE
5	AF	75	GLU
5	AF	76	THR
5	AF	78	PHE
5	AF	86	ARG
5	AF	88	MET
5	AF	92	THR
6	AG	3	ARG
6	AG	10	LYS
6	AG	11	ILE
6	AG	14	ASP
6	AG	16	LYS
6	AG	21	LEU
6	AG	22	LEU
6	AG	42	VAL
6	AG	52	ARG
6	AG	55	LYS
6	AG	72	VAL
6	AG	84	TYR
6	AG	89	GLU
6	AG	94	ARG
6	AG	112	ASP
6	AG	137	ARG
6	AG	143	MET
7	AH	26	MET
7	AH	30	LYS
7	AH	37	ASN
7	AH	39	LEU
7	AH	48	PHE
7	AH	55	LYS
7	AH	58	LEU
7	AH	82	LEU
7	AH	93	LYS
7	AH	111	THR
7	AH	113	ARG
7	AH	120	LEU
8	AI	11	ARG

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Mol	Chain	Res	Type
8	AI	26	LYS
8	AI	36	GLN
8	AI	38	PHE
8	AI	44	ARG
8	AI	56	MET
8	AI	58	GLU
8	AI	59	LYS
8	AI	84	ARG
8	AI	86	LEU
8	AI	87	MET
8	AI	106	ASP
8	AI	112	ARG
8	AI	121	ARG
8	AI	126	PHE
9	AJ	5	ARG
9	AJ	11	LYS
9	AJ	14	ASP
9	AJ	18	ILE
9	AJ	40	ILE
9	AJ	48	ARG
9	AJ	49	PHE
9	AJ	57	VAL
9	AJ	68	ARG
9	AJ	71	LEU
9	AJ	75	ASP
9	AJ	81	GLU
9	AJ	88	MET
9	AJ	89	ARG
9	AJ	92	LEU
9	AJ	102	LEU
10	AK	12	ARG
10	AK	35	ASP
10	AK	36	ARG
10	AK	37	GLN
10	AK	52	ARG
10	AK	56	LYS
10	AK	58	THR
10	AK	60	PHE
10	AK	73	VAL
10	AK	83	VAL
10	AK	84	MET
10	AK	92	ARG

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Mol	Chain	Res	Type
10	AK	99	LEU
10	AK	107	THR
10	AK	115	ILE
10	AK	117	HIS
10	AK	121	ARG
10	AK	126	ARG
11	AL	24	GLU
11	AL	28	GLN
11	AL	30	ARG
11	AL	35	ARG
11	AL	38	THR
11	AL	48	LEU
11	AL	50	LYS
11	AL	58	ASN
11	AL	71	HIS
11	AL	74	GLN
11	AL	87	LYS
11	AL	98	ARG
11	AL	102	ASP
12	AM	18	LEU
12	AM	28	ARG
12	AM	41	ASP
12	AM	43	LYS
12	AM	44	ILE
12	AM	57	ASP
12	AM	62	PHE
12	AM	67	ASP
12	AM	68	LEU
12	AM	90	HIS
12	AM	91	ARG
12	AM	97	ARG
12	AM	100	ARG
12	AM	101	THR
12	AM	102	LYS
12	AM	109	LYS
13	AN	15	LEU
13	AN	27	LYS
13	AN	41	TRP
13	AN	50	LEU
13	AN	59	GLN
13	AN	60	ARG
13	AN	65	GLN

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Mol	Chain	Res	Type
13	AN	73	LEU
13	AN	74	ARG
14	AO	7	THR
14	AO	17	ASP
14	AO	25	GLU
14	AO	53	ARG
14	AO	57	ARG
14	AO	66	LEU
14	AO	70	LYS
15	AP	4	ILE
15	AP	5	ARG
15	AP	9	HIS
15	AP	19	VAL
15	AP	26	ASN
15	AP	29	ASN
15	AP	31	ARG
15	AP	34	GLU
15	AP	46	LYS
15	AP	55	ASP
15	AP	63	GLN
16	AQ	3	LYS
16	AQ	4	ILE
16	AQ	10	ARG
16	AQ	20	ILE
16	AQ	25	GLU
16	AQ	39	ARG
16	AQ	47	ASP
16	AQ	60	ILE
16	AQ	76	ARG
16	AQ	80	LYS
17	AR	20	ILE
17	AR	23	LYS
17	AR	30	ASN
17	AR	52	ARG
18	AS	5	LYS
18	AS	14	LEU
18	AS	20	LYS
18	AS	28	LYS
18	AS	31	ARG
18	AS	35	ARG
18	AS	40	PHE
18	AS	47	THR

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Mol	Chain	Res	Type
18	AS	63	ASP
18	AS	72	GLU
18	AS	73	PHE
19	AT	4	LYS
19	AT	14	GLU
19	AT	26	MET
19	AT	29	THR
19	AT	43	LYS
19	AT	53	MET
19	AT	59	ARG
19	AT	67	HIS
19	AT	70	LYS
19	AT	85	LEU
20	AB	9	LEU
20	AB	22	TRP
20	AB	23	ASN
20	AB	31	PHE
20	AB	36	LYS
20	AB	38	HIS
20	AB	40	ILE
20	AB	46	VAL
20	AB	48	MET
20	AB	56	LEU
20	AB	57	ASN
20	AB	62	ARG
20	AB	67	LEU
20	AB	81	ASP
20	AB	90	PHE
20	AB	94	ARG
20	AB	95	TRP
20	AB	116	LEU
20	AB	121	GLN
20	AB	122	ASP
20	AB	125	PHE
20	AB	127	LYS
20	AB	128	LEU
20	AB	145	ASN
20	AB	176	ASN
20	AB	188	THR
20	AB	193	ASP
20	AB	196	ASP
20	AB	202	ASN

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Mol	Chain	Res	Type
20	AB	206	ILE
20	AB	212	TYR
20	AB	222	GLU
21	AU	11	PHE
21	AU	12	ASP
21	AU	16	ARG
21	AU	32	ARG
21	AU	41	THR
21	AU	48	LYS
21	AU	53	LYS
24	BV	2	PHE
24	BV	3	THR
24	BV	9	ARG
24	BV	35	GLU
24	BV	40	ILE
24	BV	42	LEU
24	BV	46	LYS
24	BV	51	GLN
24	BV	56	PHE
24	BV	70	ILE
24	BV	73	LYS
24	BV	79	ARG
24	BV	82	TYR
24	BV	87	GLN
25	BC	5	CYS
25	BC	14	HIS
25	BC	17	LYS
25	BC	18	VAL
25	BC	27	LYS
25	BC	28	PRO
25	BC	29	PHE
25	BC	32	LEU
25	BC	33	LEU
25	BC	35	LYS
25	BC	36	ASN
25	BC	38	LYS
25	BC	45	ASN
25	BC	51	ARG
25	BC	57	HIS
25	BC	58	LYS
25	BC	62	ARG
25	BC	64	VAL

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Mol	Chain	Res	Type
25	BC	78	GLU
25	BC	86	ARG
25	BC	94	LEU
25	BC	102	TYR
25	BC	110	LYS
25	BC	125	PRO
25	BC	127	ASN
25	BC	129	LEU
25	BC	131	MET
25	BC	144	GLU
25	BC	152	GLN
25	BC	155	ARG
25	BC	162	GLN
25	BC	171	VAL
25	BC	180	MET
25	BC	181	ARG
25	BC	186	ASP
25	BC	199	HIS
25	BC	201	LEU
25	BC	203	VAL
25	BC	206	LYS
25	BC	220	ARG
25	BC	222	THR
25	BC	224	MET
25	BC	225	ASN
25	BC	231	HIS
25	BC	235	GLU
25	BC	237	ARG
25	BC	246	PRO
25	BC	256	THR
25	BC	261	ARG
25	BC	263	ASP
25	BC	268	ARG
26	BD	18	ASP
26	BD	25	THR
26	BD	29	VAL
26	BD	30	GLU
26	BD	35	THR
26	BD	36	GLN
26	BD	38	LYS
26	BD	39	ASP
26	BD	48	ILE

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Mol	Chain	Res	Type
26	BD	52	THR
26	BD	56	LYS
26	BD	59	ARG
26	BD	62	LYS
26	BD	67	HIS
26	BD	70	LYS
26	BD	96	ILE
26	BD	103	ASP
26	BD	105	LYS
26	BD	123	LYS
26	BD	127	PHE
26	BD	128	ARG
26	BD	131	ASP
26	BD	133	THR
26	BD	138	LEU
26	BD	140	HIS
26	BD	141	ARG
26	BD	148	GLN
26	BD	155	VAL
26	BD	157	LYS
26	BD	159	LYS
26	BD	160	LYS
26	BD	167	ASN
26	BD	177	VAL
26	BD	197	THR
27	BE	2	GLU
27	BE	4	VAL
27	BE	5	LEU
27	BE	10	SER
27	BE	18	THR
27	BE	22	ASP
27	BE	24	ASN
27	BE	41	GLN
27	BE	43	THR
27	BE	44	ARG
27	BE	46	GLN
27	BE	47	LYS
27	BE	49	ARG
27	BE	57	LYS
27	BE	58	LYS
27	BE	59	PRO
27	BE	67	ARG

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Mol	Chain	Res	Type
27	BE	78	TRP
27	BE	85	PHE
27	BE	88	ARG
27	BE	91	ASP
27	BE	98	LYS
27	BE	118	LEU
27	BE	132	LYS
27	BE	134	LEU
27	BE	139	LYS
27	BE	143	LEU
27	BE	148	ILE
27	BE	162	ARG
27	BE	164	LEU
27	BE	169	VAL
27	BE	171	ASP
27	BE	184	ASP
27	BE	188	MET
27	BE	193	VAL
27	BE	196	VAL
27	BE	199	MET
27	BE	200	LEU
28	BF	5	ASP
28	BF	14	LYS
28	BF	16	MET
28	BF	21	TYR
28	BF	55	ASP
28	BF	59	ILE
28	BF	63	LYS
28	BF	76	PHE
28	BF	93	GLU
28	BF	101	ARG
28	BF	126	ASN
28	BF	132	ARG
28	BF	142	TYR
28	BF	143	ASP
28	BF	146	ASP
28	BF	147	ARG
28	BF	157	THR
28	BF	160	LYS
28	BF	166	ARG
28	BF	168	LEU
28	BF	174	PHE

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Mol	Chain	Res	Type
28	BF	177	ARG
29	BG	2	ARG
29	BG	3	VAL
29	BG	17	LYS
29	BG	18	ILE
29	BG	28	LYS
29	BG	34	ARG
29	BG	46	ASP
29	BG	49	LEU
29	BG	50	THR
29	BG	57	TYR
29	BG	59	ASP
29	BG	61	TRP
29	BG	70	LEU
29	BG	80	GLU
29	BG	94	ARG
29	BG	121	THR
29	BG	128	THR
29	BG	133	LYS
29	BG	146	ASP
29	BG	150	TYR
29	BG	154	GLU
29	BG	176	LYS
30	BH	1	MET
30	BH	5	LEU
30	BH	7	ASP
30	BH	8	LYS
30	BH	22	LYS
30	BH	28	ASN
30	BH	30	LEU
30	BH	35	LYS
30	BH	42	LYS
30	BH	47	PHE
30	BH	55	GLU
30	BH	57	LYS
30	BH	66	ASN
30	BH	68	ARG
30	BH	73	ASN
30	BH	75	LEU
30	BH	89	LYS
30	BH	112	LYS
30	BH	132	PHE

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Mol	Chain	Res	Type
31	BJ	12	LYS
31	BJ	15	TRP
31	BJ	23	LYS
31	BJ	24	THR
31	BJ	25	LEU
31	BJ	28	LEU
31	BJ	34	ARG
31	BJ	35	ARG
31	BJ	37	ARG
31	BJ	39	LYS
31	BJ	49	ASP
31	BJ	58	ASN
31	BJ	75	TYR
31	BJ	90	GLU
31	BJ	96	ARG
31	BJ	98	GLU
31	BJ	114	LEU
31	BJ	119	PHE
31	BJ	122	LEU
31	BJ	131	ASN
31	BJ	136	GLN
32	BK	9	ASN
32	BK	17	ARG
32	BK	30	ARG
32	BK	31	ARG
32	BK	32	TYR
32	BK	47	ILE
32	BK	53	LYS
32	BK	54	LYS
32	BK	58	LEU
32	BK	62	VAL
32	BK	64	ARG
32	BK	71	ARG
32	BK	79	PHE
32	BK	87	LEU
32	BK	93	GLN
32	BK	98	ARG
32	BK	108	ARG
32	BK	109	SER
32	BK	113	MET
32	BK	116	ILE
32	BK	118	LEU

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Mol	Chain	Res	Type
33	BL	2	ARG
33	BL	14	LYS
33	BL	18	ARG
33	BL	27	LEU
33	BL	29	LYS
33	BL	33	ARG
33	BL	41	ARG
33	BL	47	ARG
33	BL	51	GLU
33	BL	55	MET
33	BL	61	LEU
33	BL	62	PRO
33	BL	70	LYS
33	BL	78	ARG
33	BL	81	ASP
33	BL	82	LEU
33	BL	84	LYS
33	BL	85	VAL
33	BL	96	LYS
33	BL	103	ILE
33	BL	105	ILE
33	BL	111	ILE
33	BL	112	LEU
33	BL	118	THR
33	BL	126	ARG
33	BL	127	VAL
33	BL	132	ARG
34	BM	3	GLN
34	BM	9	PHE
34	BM	10	ARG
34	BM	11	LYS
34	BM	12	MET
34	BM	14	LYS
34	BM	17	ASN
34	BM	18	ARG
34	BM	22	GLN
34	BM	36	VAL
34	BM	38	ARG
34	BM	54	THR
34	BM	55	ARG
34	BM	63	ILE
34	BM	66	ARG

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Mol	Chain	Res	Type
34	BM	71	LYS
34	BM	86	LYS
34	BM	93	VAL
34	BM	96	ILE
34	BM	100	LYS
34	BM	105	MET
34	BM	109	PRO
34	BM	110	GLU
34	BM	117	PHE
34	BM	118	LYS
34	BM	130	PHE
34	BM	133	LYS
34	BM	135	VAL
35	BN	1	MET
35	BN	9	GLN
35	BN	10	LEU
35	BN	13	ASN
35	BN	21	PHE
35	BN	33	ILE
35	BN	34	ILE
35	BN	35	LYS
35	BN	38	LEU
35	BN	45	ARG
35	BN	47	VAL
35	BN	57	THR
35	BN	75	ILE
35	BN	76	VAL
35	BN	94	TYR
35	BN	99	LYS
35	BN	102	PHE
35	BN	112	TYR
35	BN	120	GLU
36	BO	1	MET
36	BO	3	LYS
36	BO	7	ARG
36	BO	13	ARG
36	BO	15	ARG
36	BO	16	ARG
36	BO	25	ARG
36	BO	31	THR
36	BO	32	PRO
36	BO	35	ILE

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Mol	Chain	Res	Type
36	BO	38	GLN
36	BO	40	ILE
36	BO	56	LYS
36	BO	62	LEU
36	BO	64	TYR
36	BO	69	ASP
36	BO	84	GLU
36	BO	85	LYS
36	BO	116	GLN
37	BP	2	ASN
37	BP	8	GLU
37	BP	15	ASP
37	BP	24	THR
37	BP	30	TRP
37	BP	42	PHE
37	BP	47	ILE
37	BP	49	ILE
37	BP	50	ARG
37	BP	52	ARG
37	BP	54	LEU
37	BP	55	HIS
37	BP	60	VAL
37	BP	73	PHE
37	BP	74	GLN
37	BP	87	ARG
37	BP	88	ARG
37	BP	96	LEU
37	BP	108	ARG
37	BP	113	LEU
37	BP	114	ASN
38	BQ	3	VAL
38	BQ	5	ARG
38	BQ	12	ARG
38	BQ	36	GLN
38	BQ	38	VAL
38	BQ	39	ILE
38	BQ	55	GLN
38	BQ	63	ARG
38	BQ	65	ASN
38	BQ	71	ASN
38	BQ	80	ASN
38	BQ	84	LYS

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Mol	Chain	Res	Type
38	BQ	92	LYS
38	BQ	94	LEU
38	BQ	101	ASP
38	BQ	102	LYS
38	BQ	108	LEU
39	BR	12	HIS
39	BR	13	ARG
39	BR	18	GLN
39	BR	20	VAL
39	BR	39	LEU
39	BR	40	MET
39	BR	62	GLU
39	BR	79	ARG
39	BR	83	TYR
39	BR	89	HIS
39	BR	91	GLN
39	BR	93	PHE
40	BS	6	LYS
40	BS	8	ARG
40	BS	13	SER
40	BS	28	LYS
40	BS	60	HIS
40	BS	67	ASP
40	BS	90	LYS
40	BS	94	ASP
40	BS	95	ARG
40	BS	110	ARG
41	BT	1	MET
41	BT	3	ARG
41	BT	4	GLU
41	BT	6	ARG
41	BT	8	LEU
41	BT	10	VAL
41	BT	15	HIS
41	BT	24	MET
41	BT	26	LYS
41	BT	30	ILE
41	BT	31	VAL
41	BT	50	LEU
41	BT	61	LEU
41	BT	68	LYS
41	BT	69	ARG

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Mol	Chain	Res	Type
41	BT	79	ASP
41	BT	85	VAL
41	BT	94	ASP
42	BU	3	LYS
42	BU	13	LEU
42	BU	25	LYS
42	BU	43	LYS
42	BU	52	ASN
42	BU	61	GLU
42	BU	64	ILE
42	BU	68	ASN
42	BU	71	ILE
42	BU	72	PHE
42	BU	81	ARG
42	BU	85	ARG
42	BU	87	GLU
42	BU	90	LYS
42	BU	93	ARG
42	BU	95	PHE
42	BU	98	ASN
43	BW	2	HIS
43	BW	10	ARG
43	BW	18	LYS
43	BW	23	LYS
43	BW	40	ARG
43	BW	52	CYS
43	BW	54	ARG
43	BW	65	LYS
43	BW	69	GLU
43	BW	71	LYS
43	BW	75	ASN
43	BW	81	ILE
44	BX	1	MET
44	BX	2	LYS
44	BX	8	GLU
44	BX	12	GLU
44	BX	14	LEU
44	BX	15	ASN
44	BX	17	GLU
44	BX	18	LEU
44	BX	22	LEU
44	BX	24	GLU

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Mol	Chain	Res	Type
44	BX	31	GLN
44	BX	38	GLN
44	BX	42	LEU
44	BX	43	LEU
44	BX	45	GLN
44	BX	50	VAL
44	BX	53	VAL
45	BY	6	ILE
45	BY	7	THR
45	BY	8	GLN
45	BY	9	THR
45	BY	10	ARG
45	BY	15	ARG
45	BY	16	LEU
45	BY	17	PRO
45	BY	31	ILE
45	BY	37	ARG
45	BY	39	ASP
46	BZ	6	HIS
46	BZ	8	LYS
46	BZ	11	GLU
46	BZ	13	THR
46	BZ	18	CYS
46	BZ	25	ARG
46	BZ	32	LEU
46	BZ	33	ASN
46	BZ	35	ASP
46	BZ	40	CYS
46	BZ	41	HIS
46	BZ	44	PHE
46	BZ	49	ARG
46	BZ	58	ASP
46	BZ	65	ASN
46	BZ	70	LYS
47	B0	3	GLN
47	B0	4	GLN
47	B0	5	ASN
47	B0	6	LYS
47	B0	8	THR
47	B0	16	ARG
47	B0	25	THR
47	B0	27	LEU

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Mol	Chain	Res	Type
47	B0	29	VAL
47	B0	32	THR
47	B0	52	LYS
48	B1	6	GLU
48	B1	7	LYS
48	B1	9	LYS
48	B1	20	TYR
48	B1	24	LYS
48	B1	32	LYS
48	B1	33	LEU
48	B1	34	GLU
48	B1	44	GLN
48	B1	46	VAL
48	B1	49	LYS
48	B1	52	LYS
48	B1	53	ILE
49	B2	12	ARG
49	B2	18	PHE
49	B2	19	ARG
49	B2	21	ARG
49	B2	25	LYS
49	B2	28	ARG
49	B2	29	GLN
49	B2	33	ARG
49	B2	46	LYS
50	B3	4	LYS
50	B3	6	VAL
50	B3	12	ARG
50	B3	21	PHE
50	B3	25	HIS
50	B3	28	LEU
50	B3	29	ARG
50	B3	37	THR
50	B3	41	ARG
50	B3	44	ARG
51	B4	2	LYS
51	B4	12	ARG
51	B4	18	LYS
51	B4	19	ARG
51	B4	20	ASP
51	B4	32	LYS
51	B4	33	HIS

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Mol	Chain	Res	Type
51	B4	35	GLN
51	B4	37	GLN
52	BI	63	ASP
52	BI	96	LYS
2	CC	2	GLN
2	CC	13	ILE
2	CC	19	SER
2	CC	25	THR
2	CC	28	PHE
2	CC	30	ASP
2	CC	40	GLN
2	CC	46	LEU
2	CC	87	ARG
2	CC	88	LYS
2	CC	96	VAL
2	CC	128	MET
2	CC	129	PHE
2	CC	134	LYS
2	CC	143	LEU
2	CC	164	THR
2	CC	165	GLU
2	CC	166	TRP
2	CC	168	ARG
2	CC	206	ILE
3	CD	2	ARG
3	CD	14	GLU
3	CD	25	ARG
3	CD	28	ASP
3	CD	30	LYS
3	CD	39	GLN
3	CD	46	ARG
3	CD	47	LEU
3	CD	49	ASP
3	CD	55	ARG
3	CD	58	GLN
3	CD	69	ARG
3	CD	71	PHE
3	CD	77	GLU
3	CD	84	ASN
3	CD	92	LEU
3	CD	93	LEU
3	CD	100	VAL

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Mol	Chain	Res	Type
3	CD	106	PHE
3	CD	125	ASN
3	CD	127	ARG
3	CD	130	ASN
3	CD	146	GLU
3	CD	147	LYS
3	CD	154	VAL
3	CD	160	LEU
3	CD	170	LEU
3	CD	186	GLU
3	CD	199	ILE
4	CE	9	GLU
4	CE	11	GLN
4	CE	23	THR
4	CE	25	LYS
4	CE	30	PHE
4	CE	45	VAL
4	CE	55	VAL
4	CE	61	LYS
4	CE	72	ASN
4	CE	89	THR
4	CE	92	ARG
4	CE	95	MET
4	CE	104	ILE
4	CE	120	HIS
4	CE	133	ILE
4	CE	141	ASP
4	CE	144	GLU
4	CE	146	MET
4	CE	151	MET
5	CF	2	ARG
5	CF	3	HIS
5	CF	6	ILE
5	CF	13	ASP
5	CF	39	LEU
5	CF	54	LEU
5	CF	55	HIS
5	CF	59	TYR
5	CF	76	THR
5	CF	86	ARG
5	CF	91	ARG
6	CG	2	ARG

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Mol	Chain	Res	Type
6	CG	8	GLN
6	CG	12	LEU
6	CG	46	LEU
6	CG	75	LYS
6	CG	78	ARG
6	CG	84	TYR
6	CG	89	GLU
6	CG	98	LEU
6	CG	108	ARG
6	CG	115	MET
6	CG	117	LEU
6	CG	128	GLU
6	CG	153	TYR
7	CH	8	ASP
7	CH	37	ASN
7	CH	55	LYS
7	CH	57	GLU
7	CH	68	LYS
7	CH	72	GLU
7	CH	82	LEU
7	CH	88	LYS
7	CH	107	LYS
7	CH	123	GLU
7	CH	124	ILE
8	CI	19	PHE
8	CI	26	LYS
8	CI	27	ILE
8	CI	31	GLN
8	CI	37	TYR
8	CI	45	MET
8	CI	60	LEU
8	CI	61	ASP
8	CI	62	LEU
8	CI	83	THR
8	CI	87	MET
8	CI	109	GLN
8	CI	122	ARG
8	CI	126	PHE
9	CJ	32	THR
9	CJ	37	ARG
9	CJ	47	GLU
9	CJ	71	LEU

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Mol	Chain	Res	Type
9	CJ	77	VAL
9	CJ	78	GLU
9	CJ	92	LEU
9	CJ	99	GLN
10	CK	28	ASN
10	CK	31	VAL
10	CK	34	THR
10	CK	75	GLU
10	CK	76	TYR
10	CK	78	ILE
10	CK	84	MET
10	CK	94	SER
10	CK	97	ARG
10	CK	109	ILE
10	CK	118	ASN
11	CL	8	ARG
11	CL	14	LYS
11	CL	17	LYS
11	CL	26	CYS
11	CL	28	GLN
11	CL	34	THR
11	CL	40	THR
11	CL	49	ARG
11	CL	61	GLU
11	CL	69	GLU
11	CL	75	GLU
11	CL	88	ASP
11	CL	93	ARG
11	CL	107	LYS
11	CL	111	GLN
11	CL	120	ARG
11	CL	122	LYS
12	CM	2	ARG
12	CM	15	VAL
12	CM	40	GLU
12	CM	47	LEU
12	CM	49	GLU
12	CM	56	ARG
12	CM	67	ASP
12	CM	97	ARG
12	CM	102	LYS
12	CM	109	LYS

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Mol	Chain	Res	Type
12	CM	112	ARG
13	CN	5	MET
13	CN	19	TYR
13	CN	72	PHE
13	CN	82	LYS
13	CN	89	ARG
13	CN	95	LEU
14	CO	4	THR
14	CO	39	GLN
14	CO	50	HIS
14	CO	56	LEU
14	CO	59	VAL
14	CO	63	ARG
14	CO	69	LEU
14	CO	78	THR
14	CO	83	ARG
15	CP	5	ARG
15	CP	14	ARG
15	CP	28	ARG
15	CP	32	PHE
15	CP	40	ASN
15	CP	50	THR
15	CP	57	ILE
15	CP	67	ILE
16	CQ	5	ARG
16	CQ	29	LYS
16	CQ	39	ARG
16	CQ	60	ILE
16	CQ	66	LEU
16	CQ	73	THR
16	CQ	80	LYS
17	CR	23	LYS
17	CR	25	ILE
17	CR	38	ILE
17	CR	42	ARG
17	CR	47	ARG
17	CR	51	GLN
18	CS	5	LYS
18	CS	12	LEU
18	CS	13	HIS
18	CS	20	LYS
18	CS	26	ASP

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Mol	Chain	Res	Type
18	CS	27	LYS
18	CS	28	LYS
18	CS	36	ARG
18	CS	47	THR
18	CS	52	ASN
18	CS	63	ASP
18	CS	64	GLU
18	CS	73	PHE
19	CT	9	ARG
19	CT	29	THR
19	CT	35	TYR
19	CT	51	ASN
19	CT	53	MET
19	CT	63	LYS
19	CT	69	ASN
19	CT	70	LYS
19	CT	73	ARG
19	CT	77	ASN
20	CB	8	MET
20	CB	15	PHE
20	CB	22	TRP
20	CB	23	ASN
20	CB	35	ASN
20	CB	36	LYS
20	CB	37	VAL
20	CB	46	VAL
20	CB	53	LEU
20	CB	57	ASN
20	CB	62	ARG
20	CB	84	LEU
20	CB	94	ARG
20	CB	95	TRP
20	CB	116	LEU
20	CB	121	GLN
20	CB	127	LYS
20	CB	139	GLU
20	CB	143	LEU
20	CB	145	ASN
20	CB	176	ASN
20	CB	188	THR
20	CB	191	ASP
20	CB	202	ASN

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Mol	Chain	Res	Type
20	CB	207	ARG
20	CB	213	LEU
20	CB	222	GLU
21	CU	7	GLU
21	CU	8	ASN
21	CU	17	ARG
21	CU	18	PHE
21	CU	20	ARG
21	CU	24	LYS
21	CU	33	ARG
21	CU	35	GLU
21	CU	36	PHE
21	CU	37	TYR
21	CU	38	GLU
21	CU	42	THR
21	CU	48	LYS
24	DV	18	ARG
24	DV	48	MET
24	DV	51	GLN
24	DV	69	GLU
24	DV	70	ILE
24	DV	86	LEU
25	DC	12	ARG
25	DC	28	PRO
25	DC	29	PHE
25	DC	32	LEU
25	DC	33	LEU
25	DC	34	GLU
25	DC	38	LYS
25	DC	42	ARG
25	DC	47	ARG
25	DC	49	THR
25	DC	51	ARG
25	DC	57	HIS
25	DC	58	LYS
25	DC	61	TYR
25	DC	62	ARG
25	DC	69	ASN
25	DC	70	LYS
25	DC	85	ASN
25	DC	102	TYR
25	DC	104	LEU

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Mol	Chain	Res	Type
25	DC	107	LYS
25	DC	109	LEU
25	DC	110	LYS
25	DC	125	PRO
25	DC	127	ASN
25	DC	129	LEU
25	DC	133	ASN
25	DC	141	HIS
25	DC	145	MET
25	DC	152	GLN
25	DC	155	ARG
25	DC	164	VAL
25	DC	171	VAL
25	DC	172	THR
25	DC	180	MET
25	DC	194	VAL
25	DC	201	LEU
25	DC	203	VAL
25	DC	206	LYS
25	DC	211	ARG
25	DC	231	HIS
25	DC	235	GLU
25	DC	237	ARG
25	DC	246	PRO
25	DC	250	GLN
25	DC	251	THR
25	DC	254	LYS
25	DC	259	ASN
25	DC	265	PHE
26	DD	1	MET
26	DD	4	LEU
26	DD	9	VAL
26	DD	14	ILE
26	DD	15	PHE
26	DD	27	ILE
26	DD	33	ARG
26	DD	34	VAL
26	DD	36	GLN
26	DD	40	LEU
26	DD	45	TYR
26	DD	59	ARG
26	DD	60	VAL

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Mol	Chain	Res	Type
26	DD	77	ARG
26	DD	79	LEU
26	DD	82	PHE
26	DD	92	VAL
26	DD	101	PHE
26	DD	104	VAL
26	DD	106	LYS
26	DD	114	LYS
26	DD	118	PHE
26	DD	121	THR
26	DD	124	ARG
26	DD	126	ASN
26	DD	127	PHE
26	DD	128	ARG
26	DD	133	THR
26	DD	138	LEU
26	DD	139	SER
26	DD	142	VAL
26	DD	151	THR
26	DD	152	PRO
26	DD	176	ASP
26	DD	185	ASN
26	DD	188	LEU
26	DD	202	ILE
26	DD	208	LYS
27	DE	1	MET
27	DE	19	PHE
27	DE	24	ASN
27	DE	43	THR
27	DE	49	ARG
27	DE	84	THR
27	DE	99	LYS
27	DE	100	MET
27	DE	106	LYS
27	DE	109	LEU
27	DE	115	GLN
27	DE	117	ARG
27	DE	132	LYS
27	DE	133	LEU
27	DE	134	LEU
27	DE	137	LYS
27	DE	138	LEU

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Mol	Chain	Res	Type
27	DE	143	LEU
27	DE	147	LEU
27	DE	148	ILE
27	DE	157	LEU
27	DE	158	PHE
27	DE	164	LEU
27	DE	167	VAL
27	DE	173	THR
27	DE	184	ASP
27	DE	185	LYS
27	DE	188	MET
28	DF	16	MET
28	DF	19	PHE
28	DF	32	LYS
28	DF	45	ASP
28	DF	56	LEU
28	DF	59	ILE
28	DF	76	PHE
28	DF	86	CYS
28	DF	90	LEU
28	DF	93	GLU
28	DF	105	ILE
28	DF	107	VAL
28	DF	132	ARG
28	DF	133	GLU
28	DF	135	ILE
28	DF	141	ASP
28	DF	143	ASP
28	DF	149	ARG
28	DF	161	SER
28	DF	174	PHE
28	DF	176	PHE
28	DF	177	ARG
29	DG	2	ARG
29	DG	15	ASP
29	DG	16	VAL
29	DG	17	LYS
29	DG	29	ASN
29	DG	31	GLU
29	DG	35	THR
29	DG	43	LYS
29	DG	71	LEU

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Mol	Chain	Res	Type
29	DG	72	ASN
29	DG	84	LYS
29	DG	85	LYS
29	DG	86	LEU
29	DG	94	ARG
29	DG	104	LEU
29	DG	113	ASP
29	DG	116	LEU
29	DG	121	THR
29	DG	151	ARG
29	DG	152	ARG
29	DG	154	GLU
29	DG	171	LYS
29	DG	174	LYS
30	DH	4	ILE
30	DH	7	ASP
30	DH	8	LYS
30	DH	11	ASN
30	DH	30	LEU
30	DH	38	PRO
30	DH	45	GLU
30	DH	50	ARG
30	DH	53	GLU
30	DH	57	LYS
30	DH	68	ARG
30	DH	70	GLU
30	DH	75	LEU
30	DH	96	THR
30	DH	97	ARG
30	DH	101	ASP
30	DH	112	LYS
30	DH	119	ASN
30	DH	124	THR
30	DH	129	GLU
30	DH	137	GLU
30	DH	141	LYS
31	DJ	1	MET
31	DJ	7	LYS
31	DJ	15	TRP
31	DJ	25	LEU
31	DJ	27	ARG
31	DJ	28	LEU

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Mol	Chain	Res	Type
31	DJ	31	GLU
31	DJ	35	ARG
31	DJ	36	LEU
31	DJ	41	LYS
31	DJ	47	HIS
31	DJ	61	LYS
31	DJ	67	ASN
31	DJ	69	ARG
31	DJ	85	LYS
31	DJ	89	PHE
31	DJ	91	GLU
31	DJ	95	ARG
31	DJ	109	LEU
31	DJ	118	MET
31	DJ	121	LYS
31	DJ	140	LEU
32	DK	2	ILE
32	DK	8	LEU
32	DK	10	VAL
32	DK	30	ARG
32	DK	32	TYR
32	DK	37	ASP
32	DK	58	LEU
32	DK	65	THR
32	DK	79	PHE
32	DK	98	ARG
32	DK	104	THR
32	DK	105	ARG
32	DK	120	PRO
33	DL	4	ASN
33	DL	14	LYS
33	DL	19	LEU
33	DL	21	ARG
33	DL	25	SER
33	DL	38	GLN
33	DL	39	LYS
33	DL	46	VAL
33	DL	48	ARG
33	DL	54	GLN
33	DL	55	MET
33	DL	60	ARG
33	DL	69	ARG

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Mol	Chain	Res	Type
33	DL	77	ILE
33	DL	92	LEU
33	DL	96	LYS
33	DL	104	GLN
33	DL	106	GLU
33	DL	107	PHE
33	DL	115	GLU
33	DL	119	PRO
33	DL	125	LEU
33	DL	128	THR
33	DL	135	ILE
33	DL	136	GLU
34	DM	1	MET
34	DM	5	LYS
34	DM	6	ARG
34	DM	9	PHE
34	DM	14	LYS
34	DM	16	ARG
34	DM	18	ARG
34	DM	22	GLN
34	DM	26	VAL
34	DM	28	PHE
34	DM	33	LEU
34	DM	38	ARG
34	DM	42	THR
34	DM	54	THR
34	DM	71	LYS
34	DM	72	PRO
34	DM	76	LYS
34	DM	80	VAL
34	DM	81	ARG
34	DM	84	LYS
34	DM	88	ASN
34	DM	90	GLU
34	DM	91	TYR
34	DM	96	ILE
34	DM	97	GLN
34	DM	100	LYS
34	DM	105	MET
34	DM	106	ASP
34	DM	111	GLU
34	DM	115	GLU

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Mol	Chain	Res	Type
34	DM	118	LYS
34	DM	123	LYS
34	DM	124	LEU
34	DM	127	LYS
35	DN	4	ARG
35	DN	5	LYS
35	DN	12	ARG
35	DN	18	GLN
35	DN	36	THR
35	DN	42	LYS
35	DN	49	GLU
35	DN	70	THR
35	DN	71	ARG
35	DN	79	LEU
35	DN	81	ASN
35	DN	87	PHE
35	DN	94	TYR
35	DN	96	ARG
35	DN	117	ASP
35	DN	121	LYS
36	DO	12	THR
36	DO	18	LEU
36	DO	21	LEU
36	DO	27	VAL
36	DO	30	ARG
36	DO	40	ILE
36	DO	64	TYR
36	DO	65	THR
36	DO	84	GLU
36	DO	87	ILE
36	DO	97	PHE
36	DO	98	GLN
36	DO	100	HIS
36	DO	116	GLN
37	DP	5	LYS
37	DP	8	GLU
37	DP	10	GLU
37	DP	18	SER
37	DP	19	PHE
37	DP	23	ASP
37	DP	24	THR
37	DP	27	VAL

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Mol	Chain	Res	Type
37	DP	50	ARG
37	DP	52	ARG
37	DP	54	LEU
37	DP	60	VAL
37	DP	61	ARG
37	DP	71	ARG
37	DP	73	PHE
37	DP	76	HIS
37	DP	80	VAL
37	DP	81	ASP
37	DP	84	SER
37	DP	87	ARG
37	DP	100	ARG
37	DP	101	GLU
37	DP	108	ARG
37	DP	109	ILE
37	DP	111	GLU
37	DP	112	ARG
37	DP	113	LEU
38	DQ	4	LYS
38	DQ	13	HIS
38	DQ	17	LEU
38	DQ	35	PHE
38	DQ	36	GLN
38	DQ	39	ILE
38	DQ	40	LYS
38	DQ	52	ARG
38	DQ	53	LYS
38	DQ	58	GLN
38	DQ	63	ARG
38	DQ	70	GLN
38	DQ	78	PHE
38	DQ	84	LYS
38	DQ	90	ASP
38	DQ	91	ARG
38	DQ	93	ILE
38	DQ	100	PHE
38	DQ	103	VAL
39	DR	4	VAL
39	DR	6	GLN
39	DR	7	SER
39	DR	14	VAL

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Mol	Chain	Res	Type
39	DR	18	GLN
39	DR	22	LEU
39	DR	23	GLU
39	DR	26	ASP
39	DR	32	THR
39	DR	43	ASN
39	DR	62	GLU
39	DR	75	VAL
39	DR	84	ARG
39	DR	85	LYS
39	DR	91	GLN
39	DR	99	THR
39	DR	102	SER
40	DS	7	HIS
40	DS	11	ARG
40	DS	23	LEU
40	DS	24	ILE
40	DS	25	ARG
40	DS	27	LYS
40	DS	42	LYS
40	DS	46	LEU
40	DS	73	LYS
40	DS	82	MET
40	DS	83	LYS
40	DS	85	ILE
40	DS	107	VAL
41	DT	5	GLU
41	DT	6	ARG
41	DT	11	LEU
41	DT	18	GLU
41	DT	22	THR
41	DT	24	MET
41	DT	33	LYS
41	DT	51	PHE
41	DT	64	LYS
41	DT	66	LYS
41	DT	73	ARG
41	DT	76	ARG
41	DT	88	LYS
42	DU	4	ILE
42	DU	8	ASP
42	DU	13	LEU

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Mol	Chain	Res	Type
42	DU	27	VAL
42	DU	28	LEU
42	DU	40	LEU
42	DU	52	ASN
42	DU	57	ILE
42	DU	60	LYS
42	DU	65	GLN
42	DU	71	ILE
42	DU	73	ASN
42	DU	85	ARG
42	DU	88	ASP
42	DU	90	LYS
42	DU	92	VAL
42	DU	94	PHE
42	DU	96	LYS
43	DW	2	HIS
43	DW	3	LYS
43	DW	10	ARG
43	DW	13	ARG
43	DW	18	LYS
43	DW	19	ARG
43	DW	20	LEU
43	DW	25	PHE
43	DW	37	VAL
43	DW	40	ARG
43	DW	75	ASN
44	DX	8	GLU
44	DX	17	GLU
44	DX	18	LEU
44	DX	25	GLN
44	DX	28	LEU
44	DX	30	MET
44	DX	31	GLN
44	DX	36	GLN
44	DX	39	GLN
44	DX	47	ARG
45	DY	5	LYS
45	DY	10	ARG
45	DY	13	ILE
45	DY	23	LEU
45	DY	26	LEU
45	DY	35	VAL

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Mol	Chain	Res	Type
45	DY	37	ARG
45	DY	40	THR
45	DY	43	ILE
45	DY	44	ARG
45	DY	53	MET
45	DY	57	GLU
46	DZ	10	GLU
46	DZ	20	ASN
46	DZ	24	ILE
46	DZ	28	VAL
46	DZ	31	ASP
46	DZ	34	LEU
46	DZ	50	ASP
46	DZ	64	PHE
47	D0	2	VAL
47	D0	5	ASN
47	D0	12	ARG
47	D0	15	ARG
47	D0	19	ASP
47	D0	21	LEU
47	D0	22	THR
47	D0	51	ARG
47	D0	53	VAL
48	D1	8	ILE
48	D1	18	HIS
48	D1	19	PHE
48	D1	22	THR
48	D1	24	LYS
48	D1	26	LYS
48	D1	27	ARG
48	D1	35	LEU
48	D1	44	GLN
48	D1	49	LYS
49	D2	10	LEU
49	D2	18	PHE
49	D2	19	ARG
49	D2	25	LYS
49	D2	28	ARG
49	D2	34	ARG
49	D2	39	ARG
49	D2	44	VAL
49	D2	46	LYS

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Mol	Chain	Res	Type
50	D3	6	VAL
50	D3	7	ARG
50	D3	24	LYS
50	D3	29	ARG
50	D3	32	LEU
50	D3	34	LYS
50	D3	35	LYS
50	D3	51	LYS
50	D3	53	ASP
50	D3	54	LEU
50	D3	61	LEU
50	D3	63	TYR
51	D4	1	MET
51	D4	2	LYS
51	D4	9	LYS
51	D4	17	VAL
51	D4	18	LYS
51	D4	19	ARG
51	D4	20	ASP
51	D4	23	ILE
51	D4	24	ARG
51	D4	26	ILE
51	D4	30	GLU
51	D4	31	PRO
51	D4	32	LYS
51	D4	35	GLN
51	D4	36	ARG
52	DI	2	LYS
52	DI	54	ILE
52	DI	99	LYS
52	DI	121	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (329) such sidechains are listed below:

Mol	Chain	Res	Type
2	AC	2	GLN
2	AC	5	HIS
2	AC	40	GLN
2	AC	122	GLN
2	AC	139	ASN
2	AC	184	ASN
3	AD	73	ASN

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Mol	Chain	Res	Type
3	AD	84	ASN
3	AD	88	ASN
3	AD	135	GLN
3	AD	195	ASN
4	AE	69	ASN
4	AE	72	ASN
4	AE	96	GLN
4	AE	120	HIS
4	AE	121	ASN
4	AE	131	ASN
5	AF	14	GLN
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	147	ASN
7	AH	3	GLN
8	AI	4	GLN
8	AI	24	ASN
8	AI	30	ASN
8	AI	31	GLN
8	AI	36	GLN
8	AI	49	GLN
8	AI	109	GLN
9	AJ	15	HIS
9	AJ	20	GLN
9	AJ	99	GLN
10	AK	37	GLN
10	AK	118	ASN
11	AL	28	GLN
11	AL	58	ASN
11	AL	72	ASN
13	AN	59	GLN
13	AN	61	ASN
13	AN	65	GLN
14	AO	27	GLN
14	AO	34	GLN
14	AO	36	ASN
14	AO	39	GLN
15	AP	9	HIS

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Mol	Chain	Res	Type
15	AP	18	GLN
15	AP	29	ASN
15	AP	63	GLN
17	AR	53	GLN
18	AS	55	GLN
19	AT	12	GLN
19	AT	20	ASN
19	AT	54	GLN
19	AT	60	GLN
20	AB	17	HIS
20	AB	23	ASN
20	AB	35	ASN
20	AB	41	ASN
20	AB	108	GLN
20	AB	119	GLN
20	AB	121	GLN
20	AB	145	ASN
20	AB	167	HIS
20	AB	202	ASN
24	BV	49	ASN
24	BV	51	GLN
24	BV	80	HIS
24	BV	88	HIS
25	BC	14	HIS
25	BC	36	ASN
25	BC	52	HIS
25	BC	127	ASN
25	BC	133	ASN
25	BC	162	GLN
25	BC	225	ASN
25	BC	238	ASN
25	BC	250	GLN
26	BD	32	ASN
26	BD	134	HIS
26	BD	148	GLN
26	BD	150	GLN
26	BD	164	GLN
26	BD	167	ASN
26	BD	173	GLN
26	BD	185	ASN
27	BE	156	ASN
27	BE	195	GLN

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Mol	Chain	Res	Type
28	BF	26	GLN
28	BF	36	ASN
28	BF	126	ASN
29	BG	19	ASN
29	BG	21	GLN
29	BG	29	ASN
29	BG	37	ASN
29	BG	63	GLN
29	BG	138	GLN
30	BH	2	GLN
30	BH	11	ASN
30	BH	43	ASN
30	BH	73	ASN
31	BJ	80	HIS
31	BJ	131	ASN
31	BJ	135	GLN
31	BJ	136	GLN
32	BK	3	GLN
32	BK	5	GLN
33	BL	38	GLN
33	BL	99	ASN
34	BM	17	ASN
34	BM	45	GLN
34	BM	97	GLN
35	BN	9	GLN
35	BN	73	ASN
35	BN	81	ASN
35	BN	107	ASN
36	BO	19	GLN
36	BO	38	GLN
36	BO	67	ASN
36	BO	98	GLN
36	BO	104	GLN
37	BP	2	ASN
37	BP	6	GLN
37	BP	11	GLN
37	BP	40	GLN
37	BP	114	ASN
38	BQ	36	GLN
38	BQ	43	GLN
38	BQ	51	GLN
38	BQ	55	GLN

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Mol	Chain	Res	Type
38	BQ	58	GLN
38	BQ	71	ASN
38	BQ	80	ASN
39	BR	11	GLN
40	BS	7	HIS
40	BS	9	HIS
40	BS	57	ASN
40	BS	61	ASN
41	BT	28	ASN
41	BT	48	GLN
41	BT	91	GLN
42	BU	39	ASN
42	BU	65	GLN
42	BU	68	ASN
43	BW	49	ASN
43	BW	56	HIS
44	BX	38	GLN
44	BX	41	HIS
44	BX	45	GLN
45	BY	8	GLN
45	BY	19	HIS
46	BZ	41	HIS
46	BZ	65	ASN
47	B0	3	GLN
47	B0	4	GLN
47	B0	5	ASN
48	B1	25	ASN
48	B1	45	HIS
50	B3	23	HIS
51	B4	13	ASN
51	B4	35	GLN
52	BI	11	GLN
52	BI	29	GLN
52	BI	93	ASN
2	CC	2	GLN
2	CC	31	ASN
2	CC	68	HIS
2	CC	184	ASN
2	CC	189	HIS
3	CD	39	GLN
3	CD	53	GLN
3	CD	58	GLN

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Mol	Chain	Res	Type
3	CD	99	ASN
3	CD	115	GLN
3	CD	119	HIS
3	CD	130	ASN
3	CD	151	GLN
3	CD	163	GLN
3	CD	195	ASN
4	CE	42	ASN
4	CE	72	ASN
4	CE	131	ASN
4	CE	145	ASN
5	CF	17	GLN
5	CF	46	GLN
5	CF	63	ASN
6	CG	8	GLN
6	CG	27	ASN
6	CG	67	ASN
6	CG	121	ASN
7	CH	3	GLN
7	CH	37	ASN
7	CH	75	GLN
7	CH	117	GLN
8	CI	74	GLN
8	CI	80	HIS
8	CI	109	GLN
8	CI	125	GLN
9	CJ	99	GLN
10	CK	28	ASN
10	CK	37	GLN
10	CK	63	GLN
10	CK	108	ASN
10	CK	118	ASN
11	CL	28	GLN
11	CL	45	ASN
11	CL	58	ASN
11	CL	71	HIS
12	CM	7	ASN
13	CN	34	ASN
13	CN	59	GLN
13	CN	61	ASN
14	CO	36	ASN
15	CP	26	ASN

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Mol	Chain	Res	Type
15	CP	40	ASN
15	CP	63	GLN
16	CQ	8	GLN
17	CR	51	GLN
17	CR	53	GLN
18	CS	42	ASN
18	CS	52	ASN
18	CS	55	GLN
19	CT	2	ASN
19	CT	69	ASN
19	CT	83	ASN
20	CB	23	ASN
20	CB	50	ASN
20	CB	57	ASN
20	CB	108	GLN
20	CB	119	GLN
20	CB	121	GLN
20	CB	145	ASN
20	CB	202	ASN
24	DV	51	GLN
24	DV	78	GLN
24	DV	87	GLN
25	DC	45	ASN
25	DC	59	GLN
25	DC	116	GLN
25	DC	127	ASN
25	DC	133	ASN
25	DC	141	HIS
25	DC	162	GLN
25	DC	196	ASN
25	DC	199	HIS
25	DC	225	ASN
25	DC	238	ASN
26	DD	32	ASN
26	DD	36	GLN
26	DD	94	GLN
26	DD	164	GLN
26	DD	173	GLN
27	DE	24	ASN
27	DE	30	GLN
27	DE	94	GLN
27	DE	115	GLN

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Mol	Chain	Res	Type
28	DF	36	ASN
28	DF	62	GLN
28	DF	80	GLN
28	DF	134	GLN
29	DG	29	ASN
29	DG	37	ASN
29	DG	87	GLN
29	DG	100	ASN
29	DG	138	GLN
30	DH	20	ASN
30	DH	43	ASN
30	DH	133	GLN
31	DJ	40	HIS
31	DJ	86	GLN
31	DJ	135	GLN
32	DK	3	GLN
33	DL	38	GLN
33	DL	104	GLN
34	DM	3	GLN
34	DM	17	ASN
34	DM	45	GLN
34	DM	88	ASN
34	DM	97	GLN
35	DN	3	HIS
35	DN	9	GLN
35	DN	11	ASN
35	DN	16	HIS
35	DN	18	GLN
35	DN	73	ASN
35	DN	107	ASN
36	DO	19	GLN
36	DO	67	ASN
36	DO	104	GLN
37	DP	9	GLN
37	DP	11	GLN
37	DP	55	HIS
37	DP	74	GLN
38	DQ	43	GLN
38	DQ	51	GLN
38	DQ	58	GLN
38	DQ	71	ASN
38	DQ	80	ASN

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Mol	Chain	Res	Type
39	DR	6	GLN
39	DR	43	ASN
39	DR	86	GLN
40	DS	15	GLN
40	DS	31	GLN
40	DS	40	ASN
40	DS	57	ASN
41	DT	48	GLN
42	DU	53	GLN
42	DU	65	GLN
42	DU	68	ASN
43	DW	2	HIS
43	DW	39	GLN
43	DW	45	HIS
43	DW	56	HIS
44	DX	31	GLN
44	DX	45	GLN
44	DX	58	ASN
45	DY	48	ASN
46	DZ	6	HIS
46	DZ	65	ASN
47	D0	4	GLN
48	D1	44	GLN
48	D1	45	HIS
49	D2	29	GLN
52	DI	5	GLN
52	DI	33	ASN
52	DI	93	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1529/1542 (99%)	255 (16%)	0
1	CA	1529/1542 (99%)	232 (15%)	0
22	BA	116/120 (96%)	23 (19%)	0
22	DA	116/120 (96%)	22 (18%)	0
23	BB	2837/2904 (97%)	424 (14%)	0
23	DB	2837/2904 (97%)	438 (15%)	0
All	All	8964/9132 (98%)	1394 (15%)	0

All (1394) 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	51	A
1	AA	52	C
1	AA	55	A
1	AA	61	G
1	AA	66	A
1	AA	71	A
1	AA	72	A
1	AA	75	G
1	AA	78	A
1	AA	83	C
1	AA	85	U
1	AA	86	G
1	AA	87	C
1	AA	88	U
1	AA	89	U
1	AA	95	C
1	AA	108	G
1	AA	121	U
1	AA	122	G
1	AA	131	A
1	AA	151	A
1	AA	182	A
1	AA	197	A
1	AA	209	U
1	AA	210	C
1	AA	226	G
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	257	G
1	AA	258	G

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Mol	Chain	Res	Type
1	AA	266	G
1	AA	267	C
1	AA	280	C
1	AA	289	G
1	AA	316	C
1	AA	328	C
1	AA	329	A
1	AA	330	C
1	AA	332	G
1	AA	345	C
1	AA	351	G
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	411	A
1	AA	412	A
1	AA	414	A
1	AA	415	A
1	AA	421	U
1	AA	422	C
1	AA	424	G
1	AA	429	U
1	AA	430	A
1	AA	435	A
1	AA	438	U
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

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Mol	Chain	Res	Type
1	AA	469	C
1	AA	479	U
1	AA	481	G
1	AA	484	G
1	AA	485	U
1	AA	493	A
1	AA	518	C
1	AA	522	C
1	AA	527	G
1	AA	531	U
1	AA	532	A
1	AA	547	A
1	AA	559	A
1	AA	562	U
1	AA	564	C
1	AA	572	A
1	AA	573	A
1	AA	576	C
1	AA	577	G
1	AA	639	G
1	AA	650	G
1	AA	653	U
1	AA	665	A
1	AA	687	A
1	AA	700	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	815	A
1	AA	817	C
1	AA	818	G
1	AA	819	A

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Mol	Chain	Res	Type
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	846	G
1	AA	902	G
1	AA	914	A
1	AA	926	G
1	AA	927	G
1	AA	931	C
1	AA	934	C
1	AA	935	A
1	AA	960	U
1	AA	961	U
1	AA	966	G
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	983	A
1	AA	991	U
1	AA	992	U
1	AA	993	G
1	AA	1002	G
1	AA	1004	A
1	AA	1009	U
1	AA	1022	A
1	AA	1026	G
1	AA	1030	U
1	AA	1031	C
1	AA	1034	G
1	AA	1035	A
1	AA	1045	C
1	AA	1049	U
1	AA	1050	G
1	AA	1051	C
1	AA	1054	C

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Mol	Chain	Res	Type
1	AA	1065	U
1	AA	1066	C
1	AA	1070	U
1	AA	1094	G
1	AA	1095	U
1	AA	1101	A
1	AA	1102	A
1	AA	1108	G
1	AA	1111	A
1	AA	1112	C
1	AA	1126	U
1	AA	1130	A
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	1168	U
1	AA	1179	A
1	AA	1181	G
1	AA	1183	U
1	AA	1196	A
1	AA	1197	A
1	AA	1201	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
1	AA	1226	C
1	AA	1227	A
1	AA	1228	C
1	AA	1240	U
1	AA	1241	G
1	AA	1253	G
1	AA	1256	A
1	AA	1258	G
1	AA	1261	A
1	AA	1278	G

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Mol	Chain	Res	Type
1	AA	1279	G
1	AA	1280	A
1	AA	1281	C
1	AA	1285	A
1	AA	1286	U
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	1320	C
1	AA	1322	C
1	AA	1323	G
1	AA	1336	C
1	AA	1346	A
1	AA	1347	G
1	AA	1353	G
1	AA	1362	A
1	AA	1363	A
1	AA	1364	U
1	AA	1378	C
1	AA	1379	G
1	AA	1380	U
1	AA	1381	U
1	AA	1398	A
1	AA	1404	C
1	AA	1409	C
1	AA	1419	G
1	AA	1432	G
1	AA	1446	A
1	AA	1452	C
1	AA	1493	A
1	AA	1499	A
1	AA	1503	A
1	AA	1506	U
1	AA	1507	A
1	AA	1517	G
1	AA	1519	A
1	AA	1520	C
1	AA	1529	G
1	AA	1530	G

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Mol	Chain	Res	Type
1	AA	1533	C
1	AA	1534	A
22	BA	11	C
22	BA	12	C
22	BA	13	G
22	BA	14	U
22	BA	15	A
22	BA	16	G
22	BA	25	U
22	BA	26	C
22	BA	28	C
22	BA	29	A
22	BA	30	C
22	BA	35	C
22	BA	42	C
22	BA	44	G
22	BA	45	A
22	BA	56	G
22	BA	66	A
22	BA	67	G
22	BA	84	G
22	BA	90	C
22	BA	99	A
22	BA	108	A
22	BA	109	A
23	BB	2	G
23	BB	4	U
23	BB	13	A
23	BB	14	A
23	BB	34	U
23	BB	35	G
23	BB	46	G
23	BB	49	A
23	BB	63	A
23	BB	64	A
23	BB	71	A
23	BB	74	A
23	BB	75	G
23	BB	84	A
23	BB	85	G
23	BB	91	A
23	BB	92	U

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Mol	Chain	Res	Type
23	BB	96	C
23	BB	100	U
23	BB	101	A
23	BB	102	U
23	BB	103	A
23	BB	104	A
23	BB	118	A
23	BB	119	A
23	BB	120	U
23	BB	125	A
23	BB	126	A
23	BB	140	C
23	BB	142	A
23	BB	144	A
23	BB	160	A
23	BB	162	U
23	BB	163	C
23	BB	181	A
23	BB	186	G
23	BB	196	A
23	BB	199	A
23	BB	209	C
23	BB	216	A
23	BB	221	A
23	BB	222	A
23	BB	230	G
23	BB	241	A
23	BB	248	G
23	BB	252	G
23	BB	255	A
23	BB	265	A
23	BB	266	G
23	BB	271	G
23	BB	277	G
23	BB	278	A
23	BB	279	A
23	BB	281	C
23	BB	311	A
23	BB	323	C
23	BB	329	G
23	BB	330	A
23	BB	353	C

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Mol	Chain	Res	Type
23	BB	355	U
23	BB	362	A
23	BB	371	A
23	BB	372	G
23	BB	386	G
23	BB	387	U
23	BB	395	U
23	BB	396	G
23	BB	397	U
23	BB	411	G
23	BB	412	A
23	BB	423	A
23	BB	424	G
23	BB	435	C
23	BB	444	C
23	BB	455	C
23	BB	457	A
23	BB	479	A
23	BB	480	A
23	BB	481	G
23	BB	490	C
23	BB	491	G
23	BB	498	G
23	BB	504	A
23	BB	505	A
23	BB	508	A
23	BB	509	C
23	BB	512	G
23	BB	518	G
23	BB	527	C
23	BB	531	C
23	BB	532	A
23	BB	533	G
23	BB	544	C
23	BB	545	U
23	BB	546	U
23	BB	547	A
23	BB	555	G
23	BB	563	A
23	BB	572	A
23	BB	573	U
23	BB	575	A

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Mol	Chain	Res	Type
23	BB	588	U
23	BB	603	A
23	BB	613	A
23	BB	614	A
23	BB	615	U
23	BB	616	A
23	BB	627	A
23	BB	637	A
23	BB	645	C
23	BB	646	U
23	BB	654	A
23	BB	655	A
23	BB	671	C
23	BB	686	U
23	BB	717	C
23	BB	719	C
23	BB	722	A
23	BB	730	A
23	BB	747	U
23	BB	757	G
23	BB	765	C
23	BB	775	G
23	BB	782	A
23	BB	784	G
23	BB	785	G
23	BB	805	G
23	BB	812	C
23	BB	819	A
23	BB	827	U
23	BB	828	U
23	BB	846	U
23	BB	847	U
23	BB	859	G
23	BB	872	U
23	BB	876	C
23	BB	877	A
23	BB	878	A
23	BB	899	A
23	BB	910	A
23	BB	912	C
23	BB	932	U
23	BB	933	A

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Mol	Chain	Res	Type
23	BB	934	U
23	BB	941	A
23	BB	946	C
23	BB	955	U
23	BB	961	C
23	BB	973	A
23	BB	974	G
23	BB	983	A
23	BB	985	C
23	BB	991	C
23	BB	993	G
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	1025	G
23	BB	1026	G
23	BB	1033	U
23	BB	1046	A
23	BB	1047	G
23	BB	1062	G
23	BB	1070	A
23	BB	1088	A
23	BB	1090	A
23	BB	1112	G
23	BB	1122	G
23	BB	1126	A
23	BB	1132	U
23	BB	1133	A
23	BB	1134	A
23	BB	1135	C
23	BB	1136	G
23	BB	1139	G
23	BB	1142	A
23	BB	1157	G
23	BB	1173	U
23	BB	1174	U
23	BB	1175	A
23	BB	1195	G
23	BB	1205	A

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Mol	Chain	Res	Type
23	BB	1206	G
23	BB	1211	C
23	BB	1212	G
23	BB	1218	G
23	BB	1237	A
23	BB	1247	A
23	BB	1248	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	1275	A
23	BB	1276	A
23	BB	1301	A
23	BB	1302	A
23	BB	1312	U
23	BB	1325	U
23	BB	1337	G
23	BB	1341	G
23	BB	1345	C
23	BB	1352	U
23	BB	1365	A
23	BB	1379	U
23	BB	1383	A
23	BB	1384	A
23	BB	1396	U
23	BB	1397	U
23	BB	1416	G
23	BB	1417	C
23	BB	1419	A
23	BB	1420	A
23	BB	1427	A
23	BB	1428	C
23	BB	1434	A
23	BB	1451	C
23	BB	1453	A
23	BB	1454	C
23	BB	1459	G
23	BB	1460	U
23	BB	1461	C

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Mol	Chain	Res	Type
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	1508	A
23	BB	1509	A
23	BB	1510	G
23	BB	1524	G
23	BB	1532	A
23	BB	1535	A
23	BB	1537	G
23	BB	1538	G
23	BB	1540	G
23	BB	1552	A
23	BB	1558	C
23	BB	1559	U
23	BB	1567	G
23	BB	1569	A
23	BB	1578	U
23	BB	1585	C
23	BB	1608	A
23	BB	1610	A
23	BB	1634	A
23	BB	1635	A
23	BB	1640	A
23	BB	1647	U
23	BB	1648	U
23	BB	1649	G
23	BB	1674	G
23	BB	1700	A
23	BB	1701	A
23	BB	1706	C
23	BB	1707	G
23	BB	1713	A
23	BB	1715	G
23	BB	1716	U
23	BB	1722	A
23	BB	1723	G

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Mol	Chain	Res	Type
23	BB	1729	U
23	BB	1730	C
23	BB	1731	G
23	BB	1733	G
23	BB	1738	G
23	BB	1758	U
23	BB	1764	C
23	BB	1773	A
23	BB	1781	U
23	BB	1800	C
23	BB	1816	C
23	BB	1829	A
23	BB	1833	C
23	BB	1870	C
23	BB	1873	G
23	BB	1876	A
23	BB	1884	G
23	BB	1906	G
23	BB	1914	C
23	BB	1927	A
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	1966	A
23	BB	1967	C
23	BB	1970	A
23	BB	1971	U
23	BB	1972	G
23	BB	1991	U
23	BB	1993	U
23	BB	1997	C
23	BB	2022	U
23	BB	2023	C
23	BB	2031	A
23	BB	2033	A
23	BB	2043	C
23	BB	2055	C
23	BB	2056	G

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Mol	Chain	Res	Type
23	BB	2060	A
23	BB	2061	G
23	BB	2062	A
23	BB	2069	G
23	BB	2072	C
23	BB	2102	G
23	BB	2104	C
23	BB	2105	U
23	BB	2137	U
23	BB	2140	G
23	BB	2144	G
23	BB	2145	C
23	BB	2147	A
23	BB	2149	U
23	BB	2154	A
23	BB	2181	U
23	BB	2183	A
23	BB	2187	U
23	BB	2192	U
23	BB	2198	A
23	BB	2204	G
23	BB	2212	A
23	BB	2225	A
23	BB	2238	G
23	BB	2239	G
23	BB	2259	U
23	BB	2266	A
23	BB	2273	A
23	BB	2276	G
23	BB	2278	A
23	BB	2283	C
23	BB	2286	G
23	BB	2287	A
23	BB	2297	A
23	BB	2304	G
23	BB	2305	U
23	BB	2307	G
23	BB	2318	G
23	BB	2319	G
23	BB	2320	U
23	BB	2321	U
23	BB	2322	A

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Mol	Chain	Res	Type
23	BB	2325	G
23	BB	2333	A
23	BB	2334	U
23	BB	2336	A
23	BB	2337	G
23	BB	2347	C
23	BB	2383	G
23	BB	2385	C
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	2434	A
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	2535	G
23	BB	2554	U
23	BB	2566	A
23	BB	2567	G
23	BB	2573	C
23	BB	2585	U
23	BB	2586	U
23	BB	2597	G
23	BB	2602	A
23	BB	2609	U
23	BB	2613	U
23	BB	2621	G
23	BB	2629	U
23	BB	2654	A
23	BB	2682	A

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Mol	Chain	Res	Type
23	BB	2689	U
23	BB	2690	U
23	BB	2714	G
23	BB	2725	A
23	BB	2726	A
23	BB	2744	G
23	BB	2748	A
23	BB	2751	G
23	BB	2752	C
23	BB	2757	A
23	BB	2765	A
23	BB	2778	A
23	BB	2791	G
23	BB	2796	U
23	BB	2797	U
23	BB	2800	A
23	BB	2801	G
23	BB	2808	G
23	BB	2820	A
23	BB	2821	A
23	BB	2836	U
23	BB	2850	A
23	BB	2866	U
23	BB	2867	G
23	BB	2872	A
23	BB	2873	A
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	51	A
1	CA	52	C
1	CA	55	A
1	CA	61	G
1	CA	71	A
1	CA	72	A

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Mol	Chain	Res	Type
1	CA	83	C
1	CA	87	C
1	CA	92	U
1	CA	93	U
1	CA	94	G
1	CA	108	G
1	CA	121	U
1	CA	122	G
1	CA	131	A
1	CA	151	A
1	CA	182	A
1	CA	197	A
1	CA	209	U
1	CA	210	C
1	CA	226	G
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	257	G
1	CA	258	G
1	CA	266	G
1	CA	267	C
1	CA	280	C
1	CA	289	G
1	CA	328	C
1	CA	329	A
1	CA	330	C
1	CA	332	G
1	CA	345	C
1	CA	351	G
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

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Mol	Chain	Res	Type
1	CA	397	A
1	CA	398	U
1	CA	406	G
1	CA	408	A
1	CA	409	U
1	CA	411	A
1	CA	412	A
1	CA	414	A
1	CA	415	A
1	CA	421	U
1	CA	422	C
1	CA	424	G
1	CA	429	U
1	CA	430	A
1	CA	435	A
1	CA	438	U
1	CA	461	A
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	469	C
1	CA	479	U
1	CA	481	G
1	CA	484	G
1	CA	485	U
1	CA	493	A
1	CA	518	C
1	CA	522	C
1	CA	527	G
1	CA	531	U
1	CA	532	A
1	CA	547	A
1	CA	559	A
1	CA	562	U
1	CA	564	C
1	CA	572	A
1	CA	573	A
1	CA	576	C

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Mol	Chain	Res	Type
1	CA	577	G
1	CA	639	G
1	CA	650	G
1	CA	653	U
1	CA	665	A
1	CA	687	A
1	CA	700	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
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	815	A
1	CA	817	C
1	CA	818	G
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	846	G
1	CA	907	A
1	CA	914	A
1	CA	926	G
1	CA	927	G
1	CA	934	C
1	CA	935	A
1	CA	945	G
1	CA	961	U
1	CA	969	A
1	CA	974	A

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Mol	Chain	Res	Type
1	CA	976	G
1	CA	977	A
1	CA	991	U
1	CA	992	U
1	CA	993	G
1	CA	1004	A
1	CA	1009	U
1	CA	1010	U
1	CA	1020	G
1	CA	1022	A
1	CA	1023	U
1	CA	1025	U
1	CA	1031	C
1	CA	1033	G
1	CA	1034	G
1	CA	1036	A
1	CA	1044	A
1	CA	1050	G
1	CA	1053	G
1	CA	1054	C
1	CA	1065	U
1	CA	1066	C
1	CA	1091	U
1	CA	1094	G
1	CA	1095	U
1	CA	1101	A
1	CA	1118	U
1	CA	1129	C
1	CA	1130	A
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	1144	G
1	CA	1159	U
1	CA	1160	G
1	CA	1168	U
1	CA	1178	G
1	CA	1184	G
1	CA	1196	A

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Mol	Chain	Res	Type
1	CA	1197	A
1	CA	1201	A
1	CA	1202	U
1	CA	1211	U
1	CA	1212	U
1	CA	1213	A
1	CA	1214	C
1	CA	1225	A
1	CA	1226	C
1	CA	1227	A
1	CA	1241	G
1	CA	1256	A
1	CA	1257	A
1	CA	1258	G
1	CA	1280	A
1	CA	1281	C
1	CA	1282	C
1	CA	1285	A
1	CA	1298	U
1	CA	1300	G
1	CA	1301	U
1	CA	1302	C
1	CA	1303	C
1	CA	1305	G
1	CA	1317	C
1	CA	1318	A
1	CA	1320	C
1	CA	1331	G
1	CA	1362	A
1	CA	1380	U
1	CA	1397	C
1	CA	1398	A
1	CA	1400	C
1	CA	1401	G
1	CA	1419	G
1	CA	1426	G
1	CA	1432	G
1	CA	1446	A
1	CA	1452	C
1	CA	1493	A
1	CA	1497	G
1	CA	1503	A

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Mol	Chain	Res	Type
1	CA	1505	G
1	CA	1506	U
1	CA	1507	A
1	CA	1517	G
1	CA	1520	C
1	CA	1529	G
1	CA	1530	G
1	CA	1533	C
1	CA	1534	A
22	DA	13	G
22	DA	15	A
22	DA	16	G
22	DA	18	G
22	DA	25	U
22	DA	26	C
22	DA	27	C
22	DA	28	C
22	DA	29	A
22	DA	42	C
22	DA	44	G
22	DA	52	A
22	DA	53	A
22	DA	56	G
22	DA	57	A
22	DA	66	A
22	DA	67	G
22	DA	89	U
22	DA	90	C
22	DA	91	C
22	DA	99	A
22	DA	109	A
23	DB	12	U
23	DB	27	G
23	DB	35	G
23	DB	46	G
23	DB	49	A
23	DB	51	G
23	DB	52	A
23	DB	63	A
23	DB	64	A
23	DB	71	A
23	DB	74	A

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Mol	Chain	Res	Type
23	DB	75	G
23	DB	91	A
23	DB	98	G
23	DB	99	U
23	DB	100	U
23	DB	101	A
23	DB	102	U
23	DB	103	A
23	DB	118	A
23	DB	119	A
23	DB	120	U
23	DB	124	G
23	DB	125	A
23	DB	126	A
23	DB	128	C
23	DB	139	U
23	DB	141	G
23	DB	144	A
23	DB	160	A
23	DB	162	U
23	DB	163	C
23	DB	180	G
23	DB	181	A
23	DB	196	A
23	DB	216	A
23	DB	221	A
23	DB	222	A
23	DB	223	A
23	DB	230	G
23	DB	233	A
23	DB	248	G
23	DB	250	G
23	DB	252	G
23	DB	255	A
23	DB	265	A
23	DB	266	G
23	DB	267	C
23	DB	271	G
23	DB	277	G
23	DB	278	A
23	DB	289	G
23	DB	299	A

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Mol	Chain	Res	Type
23	DB	301	G
23	DB	302	C
23	DB	311	A
23	DB	312	G
23	DB	323	C
23	DB	329	G
23	DB	330	A
23	DB	333	G
23	DB	346	A
23	DB	353	C
23	DB	354	A
23	DB	361	G
23	DB	363	G
23	DB	365	U
23	DB	371	A
23	DB	372	G
23	DB	386	G
23	DB	387	U
23	DB	396	G
23	DB	403	U
23	DB	404	A
23	DB	405	U
23	DB	406	G
23	DB	411	G
23	DB	412	A
23	DB	424	G
23	DB	444	C
23	DB	448	U
23	DB	449	A
23	DB	455	C
23	DB	456	C
23	DB	457	A
23	DB	458	G
23	DB	475	C
23	DB	479	A
23	DB	480	A
23	DB	481	G
23	DB	491	G
23	DB	504	A
23	DB	505	A
23	DB	506	G
23	DB	508	A

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Mol	Chain	Res	Type
23	DB	509	C
23	DB	512	G
23	DB	527	C
23	DB	530	G
23	DB	531	C
23	DB	532	A
23	DB	533	G
23	DB	544	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	572	A
23	DB	573	U
23	DB	575	A
23	DB	588	U
23	DB	603	A
23	DB	613	A
23	DB	614	A
23	DB	615	U
23	DB	616	A
23	DB	627	A
23	DB	637	A
23	DB	645	C
23	DB	646	U
23	DB	654	A
23	DB	655	A
23	DB	671	C
23	DB	686	U
23	DB	704	G
23	DB	718	A
23	DB	719	C
23	DB	727	A
23	DB	730	A
23	DB	747	U
23	DB	757	G
23	DB	765	C
23	DB	775	G
23	DB	782	A

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Mol	Chain	Res	Type
23	DB	784	G
23	DB	785	G
23	DB	805	G
23	DB	812	C
23	DB	819	A
23	DB	827	U
23	DB	828	U
23	DB	844	A
23	DB	846	U
23	DB	847	U
23	DB	859	G
23	DB	869	G
23	DB	877	A
23	DB	899	A
23	DB	910	A
23	DB	912	C
23	DB	919	U
23	DB	931	U
23	DB	932	U
23	DB	933	A
23	DB	941	A
23	DB	946	C
23	DB	955	U
23	DB	961	C
23	DB	973	A
23	DB	974	G
23	DB	982	C
23	DB	983	A
23	DB	991	C
23	DB	995	C
23	DB	1012	U
23	DB	1013	C
23	DB	1022	G
23	DB	1025	G
23	DB	1033	U
23	DB	1046	A
23	DB	1047	G
23	DB	1070	A
23	DB	1088	A
23	DB	1090	A
23	DB	1098	A
23	DB	1112	G

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Mol	Chain	Res	Type
23	DB	1116	G
23	DB	1130	U
23	DB	1132	U
23	DB	1133	A
23	DB	1134	A
23	DB	1136	G
23	DB	1142	A
23	DB	1176	U
23	DB	1179	G
23	DB	1195	G
23	DB	1205	A
23	DB	1206	G
23	DB	1210	G
23	DB	1211	C
23	DB	1212	G
23	DB	1225	G
23	DB	1237	A
23	DB	1241	A
23	DB	1242	U
23	DB	1247	A
23	DB	1248	G
23	DB	1249	U
23	DB	1250	G
23	DB	1251	C
23	DB	1252	G
23	DB	1253	A
23	DB	1256	G
23	DB	1266	G
23	DB	1271	G
23	DB	1272	A
23	DB	1275	A
23	DB	1276	A
23	DB	1301	A
23	DB	1302	A
23	DB	1312	U
23	DB	1325	U
23	DB	1337	G
23	DB	1341	G
23	DB	1345	C
23	DB	1352	U
23	DB	1365	A
23	DB	1379	U

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Mol	Chain	Res	Type
23	DB	1383	A
23	DB	1384	A
23	DB	1396	U
23	DB	1397	U
23	DB	1416	G
23	DB	1417	C
23	DB	1419	A
23	DB	1420	A
23	DB	1427	A
23	DB	1428	C
23	DB	1434	A
23	DB	1451	C
23	DB	1453	A
23	DB	1454	C
23	DB	1455	G
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	1508	A
23	DB	1509	A
23	DB	1510	G
23	DB	1524	G
23	DB	1532	A
23	DB	1535	A
23	DB	1537	G
23	DB	1538	G
23	DB	1540	G
23	DB	1552	A
23	DB	1558	C
23	DB	1559	U
23	DB	1567	G
23	DB	1569	A
23	DB	1578	U
23	DB	1585	C

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Mol	Chain	Res	Type
23	DB	1608	A
23	DB	1610	A
23	DB	1613	G
23	DB	1634	A
23	DB	1635	A
23	DB	1640	A
23	DB	1647	U
23	DB	1648	U
23	DB	1649	G
23	DB	1674	G
23	DB	1700	A
23	DB	1706	C
23	DB	1707	G
23	DB	1713	A
23	DB	1715	G
23	DB	1716	U
23	DB	1723	G
23	DB	1729	U
23	DB	1730	C
23	DB	1731	G
23	DB	1733	G
23	DB	1738	G
23	DB	1758	U
23	DB	1764	C
23	DB	1773	A
23	DB	1781	U
23	DB	1786	A
23	DB	1800	C
23	DB	1816	C
23	DB	1829	A
23	DB	1833	C
23	DB	1870	C
23	DB	1873	G
23	DB	1876	A
23	DB	1884	G
23	DB	1906	G
23	DB	1913	A
23	DB	1914	C
23	DB	1927	A
23	DB	1929	G
23	DB	1930	G
23	DB	1937	A

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Mol	Chain	Res	Type
23	DB	1938	A
23	DB	1939	U
23	DB	1940	U
23	DB	1955	U
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	1993	U
23	DB	1997	C
23	DB	2022	U
23	DB	2023	C
23	DB	2031	A
23	DB	2033	A
23	DB	2043	C
23	DB	2055	C
23	DB	2056	G
23	DB	2060	A
23	DB	2061	G
23	DB	2062	A
23	DB	2069	G
23	DB	2076	U
23	DB	2077	A
23	DB	2092	U
23	DB	2094	A
23	DB	2096	C
23	DB	2100	G
23	DB	2104	C
23	DB	2106	U
23	DB	2107	G
23	DB	2109	U
23	DB	2134	A
23	DB	2135	A
23	DB	2138	G
23	DB	2143	C
23	DB	2145	C
23	DB	2147	A
23	DB	2156	G
23	DB	2157	G
23	DB	2181	U

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Mol	Chain	Res	Type
23	DB	2183	A
23	DB	2190	G
23	DB	2199	A
23	DB	2204	G
23	DB	2210	U
23	DB	2212	A
23	DB	2225	A
23	DB	2239	G
23	DB	2253	G
23	DB	2268	A
23	DB	2283	C
23	DB	2287	A
23	DB	2288	A
23	DB	2297	A
23	DB	2305	U
23	DB	2307	G
23	DB	2308	G
23	DB	2309	A
23	DB	2310	C
23	DB	2319	G
23	DB	2321	U
23	DB	2322	A
23	DB	2325	G
23	DB	2333	A
23	DB	2334	U
23	DB	2336	A
23	DB	2337	G
23	DB	2347	C
23	DB	2383	G
23	DB	2385	C
23	DB	2402	U
23	DB	2403	C
23	DB	2406	A
23	DB	2423	U
23	DB	2426	A
23	DB	2429	G
23	DB	2430	A
23	DB	2434	A
23	DB	2441	U
23	DB	2448	A
23	DB	2472	G
23	DB	2476	A

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Mol	Chain	Res	Type
23	DB	2491	U
23	DB	2492	U
23	DB	2502	G
23	DB	2505	G
23	DB	2506	U
23	DB	2518	A
23	DB	2529	G
23	DB	2535	G
23	DB	2554	U
23	DB	2566	A
23	DB	2567	G
23	DB	2573	C
23	DB	2585	U
23	DB	2586	U
23	DB	2597	G
23	DB	2602	A
23	DB	2609	U
23	DB	2610	C
23	DB	2613	U
23	DB	2621	G
23	DB	2629	U
23	DB	2654	A
23	DB	2682	A
23	DB	2689	U
23	DB	2690	U
23	DB	2714	G
23	DB	2725	A
23	DB	2726	A
23	DB	2744	G
23	DB	2757	A
23	DB	2765	A
23	DB	2778	A
23	DB	2791	G
23	DB	2796	U
23	DB	2797	U
23	DB	2800	A
23	DB	2801	G
23	DB	2808	G
23	DB	2820	A
23	DB	2821	A
23	DB	2836	U
23	DB	2850	A

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Mol	Chain	Res	Type
23	DB	2866	U
23	DB	2867	G
23	DB	2872	A
23	DB	2873	A
23	DB	2883	A
23	DB	2894	G
23	DB	2903	U

There are no RNA pucker outliers to report.

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 345 ligands modelled in this entry, 343 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	$\# Z > 2$
53	KSG	AA	1601	-	25,27,27	2.50	5 (20%)	30,40,40	1.09	3 (10%)
53	KSG	CA	1601	-	25,27,27	2.36	5 (20%)	30,40,40	0.99	1 (3%)

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
53	KSG	AA	1601	-	-	1/8/52/52	0/2/2/2
53	KSG	CA	1601	-	-	1/8/52/52	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
53	AA	1601	KSG	C14-C13	-9.97	1.49	1.53
53	CA	1601	KSG	C14-C13	-8.54	1.49	1.53
53	AA	1601	KSG	C13-N2	-2.94	1.31	1.36
53	CA	1601	KSG	C13-N2	-2.55	1.31	1.36
53	AA	1601	KSG	C4-C3	2.07	1.57	1.52
53	AA	1601	KSG	O7-C1	2.29	1.47	1.41
53	CA	1601	KSG	O1-C1	2.47	1.48	1.41
53	CA	1601	KSG	C1-C12	2.56	1.57	1.52
53	AA	1601	KSG	C1-C12	2.62	1.57	1.52
53	CA	1601	KSG	O7-C1	2.65	1.48	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
53	AA	1601	KSG	C1-O1-C2	-2.27	112.46	118.00
53	AA	1601	KSG	C5-C4-C3	2.06	114.47	110.84
53	CA	1601	KSG	C6-C7-C2	2.25	114.26	109.61
53	AA	1601	KSG	C4-C3-C2	2.45	114.68	109.61

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
53	AA	1601	KSG	C2-O1-C1-C12
53	CA	1601	KSG	C2-O1-C1-C12

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
53	AA	1601	KSG	3	0
53	CA	1601	KSG	1	0

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%)	0.23	23 (1%) 74 68	15, 73, 149, 180	0
1	CA	1530/1542 (99%)	0.01	14 (0%) 84 78	7, 52, 134, 180	0
2	AC	206/233 (88%)	0.83	26 (12%) 4 6	7, 68, 138, 180	0
2	CC	206/233 (88%)	0.50	11 (5%) 27 24	6, 74, 126, 169	0
3	AD	205/206 (99%)	1.06	46 (22%) 1 1	27, 79, 147, 180	0
3	CD	205/206 (99%)	0.28	1 (0%) 90 87	5, 55, 113, 165	0
4	AE	150/167 (89%)	0.61	8 (5%) 27 24	15, 62, 120, 162	0
4	CE	150/167 (89%)	0.57	8 (5%) 27 24	6, 55, 112, 180	0
5	AF	100/135 (74%)	0.59	6 (6%) 23 21	18, 72, 133, 166	0
5	CF	100/135 (74%)	0.85	13 (13%) 4 6	12, 71, 132, 165	0
6	AG	150/179 (83%)	0.80	27 (18%) 2 2	22, 89, 143, 180	0
6	CG	152/179 (84%)	0.82	21 (13%) 3 5	19, 86, 147, 180	0
7	AH	129/130 (99%)	1.19	27 (20%) 1 1	14, 70, 141, 176	0
7	CH	129/130 (99%)	0.67	18 (13%) 3 5	5, 54, 114, 160	0
8	AI	127/130 (97%)	1.11	26 (20%) 1 1	18, 90, 146, 180	0
8	CI	127/130 (97%)	1.16	29 (22%) 1 1	23, 92, 156, 180	0
9	AJ	98/103 (95%)	1.23	26 (26%) 1 1	20, 82, 157, 180	0
9	CJ	98/103 (95%)	0.66	5 (5%) 29 25	34, 79, 135, 151	0
10	AK	117/129 (90%)	0.56	7 (5%) 23 21	7, 55, 138, 160	0
10	CK	117/129 (90%)	0.42	6 (5%) 29 25	11, 50, 105, 142	0
11	AL	123/124 (99%)	0.70	10 (8%) 13 14	21, 61, 128, 178	0
11	CL	123/124 (99%)	0.83	11 (8%) 10 12	5, 39, 119, 160	0
12	AM	114/118 (96%)	1.08	19 (16%) 2 2	44, 106, 159, 173	0
12	CM	113/118 (95%)	1.26	22 (19%) 1 2	26, 100, 157, 180	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
13	AN	96/101 (95%)	1.76	34 (35%)	0	0	15, 84, 151, 180	0
13	CN	96/101 (95%)	1.10	16 (16%)	2	2	21, 82, 139, 161	0
14	AO	88/89 (98%)	0.97	17 (19%)	1	2	18, 70, 114, 180	0
14	CO	88/89 (98%)	0.53	4 (4%)	34	29	14, 51, 112, 150	0
15	AP	82/82 (100%)	0.88	11 (13%)	4	5	27, 89, 142, 180	0
15	CP	80/82 (97%)	0.93	10 (12%)	4	6	5, 46, 145, 180	0
16	AQ	80/84 (95%)	1.47	19 (23%)	1	1	37, 91, 140, 180	0
16	CQ	81/84 (96%)	0.89	7 (8%)	11	12	12, 57, 119, 154	0
17	AR	55/75 (73%)	1.30	14 (25%)	1	1	18, 69, 147, 164	0
17	CR	55/75 (73%)	0.99	7 (12%)	4	6	13, 53, 126, 175	0
18	AS	79/92 (85%)	2.06	35 (44%)	0	0	59, 112, 174, 180	0
18	CS	80/92 (86%)	2.03	32 (40%)	0	0	48, 109, 172, 180	0
19	AT	85/87 (97%)	1.12	17 (20%)	1	1	36, 94, 142, 170	0
19	CT	85/87 (97%)	1.51	25 (29%)	1	1	17, 55, 118, 179	0
20	AB	218/241 (90%)	0.86	35 (16%)	2	3	28, 92, 142, 180	0
20	CB	218/241 (90%)	0.63	21 (9%)	9	10	27, 96, 157, 180	0
21	AU	51/71 (71%)	1.08	10 (19%)	1	2	26, 94, 153, 174	0
21	CU	51/71 (71%)	1.26	11 (21%)	1	1	29, 82, 163, 180	0
22	BA	117/120 (97%)	0.04	2 (1%)	70	64	39, 67, 125, 171	0
22	DA	117/120 (97%)	-0.01	2 (1%)	70	64	34, 73, 116, 180	0
23	BB	2841/2904 (97%)	0.17	41 (1%)	75	70	9, 53, 144, 180	0
23	DB	2841/2904 (97%)	0.12	21 (0%)	87	83	5, 45, 143, 180	0
24	BV	94/94 (100%)	0.70	13 (13%)	3	5	26, 78, 134, 155	0
24	DV	94/94 (100%)	0.52	5 (5%)	27	24	34, 83, 139, 162	0
25	BC	267/273 (97%)	1.20	50 (18%)	1	2	5, 53, 143, 180	0
25	DC	267/273 (97%)	1.10	43 (16%)	2	3	5, 52, 158, 180	0
26	BD	209/209 (100%)	1.43	55 (26%)	1	1	9, 82, 157, 180	0
26	DD	209/209 (100%)	1.24	46 (22%)	1	1	5, 57, 142, 180	0
27	BE	201/201 (100%)	1.34	46 (22%)	1	1	5, 85, 162, 180	0
27	DE	201/201 (100%)	1.36	43 (21%)	1	1	6, 82, 170, 180	0
28	BF	178/179 (99%)	0.92	36 (20%)	1	1	39, 105, 168, 180	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	DF	178/179 (99%)	0.77	17 (9%) 9 10	29, 96, 154, 180	0
29	BG	176/177 (99%)	1.30	41 (23%) 1 1	34, 100, 164, 180	0
29	DG	176/177 (99%)	0.90	29 (16%) 2 3	23, 88, 156, 180	0
30	BH	149/149 (100%)	2.37	66 (44%) 0 0	22, 126, 180, 180	0
30	DH	149/149 (100%)	1.53	47 (31%) 0 1	30, 106, 167, 180	0
31	BJ	140/142 (98%)	1.01	28 (20%) 1 1	11, 88, 153, 180	0
31	DJ	140/142 (98%)	0.85	20 (14%) 3 4	7, 66, 148, 176	0
32	BK	121/123 (98%)	0.73	12 (9%) 8 9	13, 60, 118, 152	0
32	DK	121/123 (98%)	0.57	3 (2%) 58 51	5, 36, 92, 134	0
33	BL	144/144 (100%)	2.61	67 (46%) 0 0	15, 86, 167, 180	0
33	DL	144/144 (100%)	2.14	55 (38%) 0 0	5, 81, 156, 180	0
34	BM	136/136 (100%)	1.57	39 (28%) 1 1	13, 71, 175, 180	0
34	DM	136/136 (100%)	1.16	22 (16%) 2 3	11, 70, 164, 180	0
35	BN	127/127 (100%)	1.55	41 (32%) 0 0	17, 78, 169, 180	0
35	DN	127/127 (100%)	1.07	16 (12%) 4 6	5, 48, 159, 180	0
36	BO	117/117 (100%)	1.81	35 (29%) 1 1	26, 91, 167, 180	0
36	DO	117/117 (100%)	1.79	37 (31%) 0 1	5, 91, 172, 180	0
37	BP	114/115 (99%)	1.84	47 (41%) 0 0	18, 98, 172, 180	0
37	DP	114/115 (99%)	1.49	30 (26%) 1 1	7, 65, 166, 180	0
38	BQ	117/118 (99%)	0.53	10 (8%) 11 13	13, 66, 146, 174	0
38	DQ	117/118 (99%)	0.90	17 (14%) 3 4	6, 57, 158, 180	0
39	BR	103/103 (100%)	1.11	21 (20%) 1 1	24, 105, 167, 180	0
39	DR	103/103 (100%)	1.69	35 (33%) 0 0	25, 92, 157, 180	0
40	BS	110/110 (100%)	0.67	8 (7%) 16 16	9, 63, 133, 180	0
40	DS	110/110 (100%)	0.77	8 (7%) 16 16	5, 46, 137, 180	0
41	BT	99/100 (99%)	1.66	36 (36%) 0 0	25, 81, 162, 180	0
41	DT	99/100 (99%)	2.62	56 (56%) 0 0	15, 86, 168, 180	0
42	BU	102/104 (98%)	2.28	39 (38%) 0 0	17, 91, 169, 180	0
42	DU	102/104 (98%)	1.84	33 (32%) 0 0	30, 110, 171, 180	0
43	BW	84/85 (98%)	2.67	38 (45%) 0 0	10, 86, 159, 180	0
43	DW	84/85 (98%)	2.44	32 (38%) 0 0	16, 86, 168, 180	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
44	BX	63/63 (100%)	1.79	19 (30%)	1	1	20, 97, 160, 180	0
44	DX	63/63 (100%)	2.73	26 (41%)	0	0	38, 110, 169, 180	0
45	BY	58/59 (98%)	1.60	17 (29%)	1	1	13, 86, 157, 180	0
45	DY	58/59 (98%)	1.23	12 (20%)	1	1	5, 79, 136, 150	0
46	BZ	70/70 (100%)	0.91	10 (14%)	3	4	19, 66, 145, 180	0
46	DZ	70/70 (100%)	1.43	23 (32%)	0	0	5, 68, 141, 180	0
47	B0	56/57 (98%)	1.42	13 (23%)	1	1	26, 95, 180, 180	0
47	D0	56/57 (98%)	1.62	14 (25%)	1	1	10, 70, 159, 177	0
48	B1	54/55 (98%)	1.45	12 (22%)	1	1	24, 97, 160, 180	0
48	D1	54/55 (98%)	2.76	29 (53%)	0	0	8, 82, 158, 180	0
49	B2	46/46 (100%)	1.36	11 (23%)	1	1	6, 51, 154, 174	0
49	D2	46/46 (100%)	1.14	6 (13%)	4	6	9, 55, 131, 150	0
50	B3	64/65 (98%)	1.99	22 (34%)	0	0	19, 61, 155, 180	0
50	D3	64/65 (98%)	1.98	23 (35%)	0	0	5, 61, 139, 178	0
51	B4	38/38 (100%)	2.78	21 (55%)	0	0	30, 102, 170, 180	0
51	D4	38/38 (100%)	2.71	17 (44%)	0	0	14, 111, 175, 180	0
52	BI	141/142 (99%)	2.55	85 (60%)	0	0	52, 152, 180, 180	0
52	DI	141/142 (99%)	2.74	94 (66%)	0	0	83, 160, 180, 180	0
All	All	20439/21086 (96%)	0.75	2577 (12%)	4	6	5, 67, 155, 180	0

All (2577) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
33	BL	98	ALA	17.1
36	BO	58	ILE	16.8
33	BL	8	PRO	15.8
41	DT	4	GLU	15.0
42	BU	29	SER	14.8
42	BU	1	ALA	14.3
33	DL	11	GLY	14.2
44	DX	31	GLN	14.1
27	DE	171	ASP	13.1
36	BO	59	ALA	13.1
44	DX	36	GLN	12.9
36	DO	59	ALA	12.7
25	BC	5	CYS	12.7

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Mol	Chain	Res	Type	RSRZ
36	DO	58	ILE	12.7
45	BY	3	THR	12.4
42	DU	1	ALA	12.1
33	BL	7	SER	12.0
42	DU	61	GLU	11.9
48	D1	4	ILE	11.8
36	DO	60	GLU	11.7
43	BW	52	CYS	11.7
33	BL	99	ASN	11.5
42	BU	89	GLY	11.3
43	DW	52	CYS	11.2
43	BW	45	HIS	10.9
33	BL	9	ALA	10.9
45	DY	1	ALA	10.9
37	BP	24	THR	10.9
36	BO	69	ASP	10.7
33	BL	1	MET	10.7
30	BH	93	SER	10.6
43	BW	83	ALA	10.5
25	BC	7	PRO	10.4
44	DX	34	SER	10.4
44	DX	32	ALA	10.4
34	DM	133	LYS	10.1
27	BE	2	GLU	10.0
27	BE	1	MET	9.9
39	DR	44	GLY	9.8
30	BH	80	ILE	9.7
44	DX	33	ALA	9.6
37	DP	1	SER	9.6
33	BL	82	LEU	9.6
33	DL	1	MET	9.6
35	DN	124	ALA	9.5
27	DE	170	ARG	9.4
30	BH	92	GLY	9.4
42	DU	27	VAL	9.3
34	BM	56	ALA	9.3
50	B3	63	TYR	9.2
41	DT	2	ILE	9.1
33	DL	99	ASN	9.1
33	DL	4	ASN	9.0
26	DD	70	LYS	9.0
33	BL	83	ALA	9.0

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Mol	Chain	Res	Type	RSRZ
37	DP	97	TYR	9.0
27	DE	3	LEU	9.0
30	BH	147	VAL	9.0
51	B4	23	ILE	8.9
35	BN	67	PHE	8.9
28	DF	74	ALA	8.8
51	D4	31	PRO	8.8
25	DC	7	PRO	8.8
51	D4	1	MET	8.8
30	BH	124	THR	8.7
35	BN	125	ALA	8.7
23	BB	1175	A	8.6
27	DE	169	VAL	8.6
44	DX	1	MET	8.6
35	DN	126	ALA	8.5
23	BB	139	U	8.5
30	BH	91	PHE	8.4
50	B3	61	LEU	8.4
42	DU	29	SER	8.4
41	DT	70	HIS	8.4
43	DW	58	LEU	8.4
42	BU	88	ASP	8.4
43	DW	83	ALA	8.3
44	DX	20	ASN	8.3
44	DX	37	LEU	8.3
25	BC	166	ARG	8.3
35	BN	124	ALA	8.3
44	DX	35	GLY	8.3
37	BP	26	GLU	8.3
48	D1	5	ARG	8.2
43	BW	2	HIS	8.2
33	DL	6	LEU	8.2
33	DL	12	SER	8.2
26	BD	91	THR	8.2
44	BX	1	MET	8.2
42	BU	2	ALA	8.2
23	DB	2133	G	8.1
33	BL	97	ALA	8.1
33	DL	55	MET	8.1
30	BH	94	ILE	8.1
26	DD	51	THR	8.1
7	AH	1	SER	8.1

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Mol	Chain	Res	Type	RSRZ
30	DH	20	ASN	8.0
33	BL	77	ILE	8.0
43	BW	78	PHE	8.0
50	B3	64	ALA	8.0
50	B3	60	CYS	8.0
43	BW	79	ILE	8.0
13	CN	35	ALA	8.0
30	BH	85	GLY	8.0
50	D3	63	TYR	7.9
48	D1	53	ILE	7.9
33	DL	7	SER	7.9
48	D1	3	GLY	7.9
29	BG	58	ALA	7.9
27	DE	145	ASP	7.8
33	DL	144	GLU	7.8
51	B4	29	ALA	7.8
26	DD	90	PHE	7.8
35	BN	127	GLU	7.8
47	D0	55	ALA	7.8
48	B1	3	GLY	7.8
50	D3	64	ALA	7.7
33	DL	100	ILE	7.7
43	BW	82	GLU	7.7
26	DD	71	ALA	7.7
30	BH	142	VAL	7.6
30	BH	130	VAL	7.6
39	DR	26	ASP	7.6
26	BD	92	VAL	7.6
42	BU	30	SER	7.5
34	BM	1	MET	7.5
43	BW	58	LEU	7.5
36	DO	1	MET	7.5
42	BU	61	GLU	7.5
34	BM	136	MET	7.4
37	DP	112	ARG	7.4
8	CI	129	ARG	7.4
27	DE	2	GLU	7.3
52	DI	1	ALA	7.3
41	DT	93	LEU	7.3
51	B4	38	GLY	7.3
39	DR	43	ASN	7.3
43	DW	35	ILE	7.2

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Mol	Chain	Res	Type	RSRZ
47	D0	56	LYS	7.2
23	BB	2145	C	7.2
43	DW	5	ALA	7.2
36	BO	57	ALA	7.2
43	BW	1	ALA	7.2
27	BE	144	GLU	7.2
25	BC	21	PRO	7.1
25	DC	8	THR	7.1
33	DL	5	THR	7.1
42	BU	59	GLU	7.1
47	D0	33	SER	7.1
48	D1	2	LYS	7.1
43	BW	4	LYS	7.1
41	DT	1	MET	7.1
52	DI	98	GLY	7.1
34	BM	134	THR	7.1
37	DP	59	THR	7.1
25	DC	5	CYS	7.0
33	DL	74	THR	7.0
48	D1	1	ALA	7.0
33	BL	5	THR	7.0
26	DD	91	THR	7.0
35	DN	120	GLU	7.0
16	AQ	5	ARG	7.0
15	AP	81	ALA	6.9
51	D4	30	GLU	6.9
39	DR	53	PHE	6.9
6	AG	76	SER	6.9
26	BD	134	HIS	6.9
36	BO	6	ALA	6.9
42	DU	60	LYS	6.8
34	DM	10	ARG	6.8
39	DR	78	ARG	6.8
48	B1	1	ALA	6.8
34	BM	135	VAL	6.8
14	AO	88	ARG	6.7
27	BE	201	ALA	6.7
44	BX	63	ALA	6.7
23	BB	137	U	6.7
30	BH	81	ALA	6.7
42	BU	28	LEU	6.7
51	B4	1	MET	6.7

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Mol	Chain	Res	Type	RSRZ
33	BL	56	PRO	6.7
30	BH	148	ALA	6.7
29	BG	1	SER	6.6
26	BD	135	GLY	6.6
42	DU	28	LEU	6.6
49	B2	15	SER	6.6
26	BD	90	PHE	6.6
33	DL	52	GLY	6.6
34	BM	132	THR	6.6
36	BO	7	ARG	6.6
52	BI	1	ALA	6.6
18	AS	73	PHE	6.6
47	D0	44	ALA	6.6
25	BC	242	HIS	6.6
34	DM	134	THR	6.6
30	BH	86	ASP	6.5
33	BL	28	GLY	6.5
25	BC	6	LYS	6.5
12	CM	3	ILE	6.5
41	DT	71	GLY	6.5
8	AI	34	LEU	6.5
13	AN	60	ARG	6.5
48	B1	4	ILE	6.5
27	BE	114	ARG	6.5
22	DA	88	C	6.5
35	BN	122	ALA	6.5
1	AA	86	G	6.4
25	DC	268	ARG	6.4
43	DW	84	GLU	6.4
43	BW	59	PHE	6.4
23	BB	2308	G	6.4
23	DB	139	U	6.4
6	AG	78	ARG	6.4
18	CS	25	GLY	6.4
30	DH	122	LEU	6.4
35	BN	126	ALA	6.3
41	DT	99	ALA	6.3
52	BI	59	THR	6.3
43	DW	1	ALA	6.3
43	DW	68	PHE	6.3
51	D4	33	HIS	6.3
52	DI	137	LEU	6.3

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Mol	Chain	Res	Type	RSRZ
23	DB	1175	A	6.3
52	BI	58	ILE	6.2
25	BC	268	ARG	6.2
44	DX	46	VAL	6.2
37	DP	26	GLU	6.2
25	DC	6	LYS	6.2
12	CM	47	LEU	6.2
18	AS	40	PHE	6.2
48	D1	54	LYS	6.2
33	BL	49	GLY	6.2
30	BH	79	THR	6.2
44	BX	38	GLN	6.2
43	DW	73	PRO	6.1
52	DI	83	ALA	6.1
35	DN	119	SER	6.1
46	DZ	58	ASP	6.1
23	DB	645	C	6.1
44	DX	38	GLN	6.0
25	BC	8	THR	6.0
27	DE	19	PHE	6.0
18	CS	26	ASP	6.0
35	DN	123	GLU	6.0
36	BO	88	LYS	5.9
33	BL	38	GLN	5.9
49	B2	14	ARG	5.9
37	DP	4	ILE	5.9
17	CR	19	GLU	5.9
35	DN	127	GLU	5.9
29	DG	43	LYS	5.9
27	BE	42	GLY	5.9
25	BC	256	THR	5.9
26	DD	69	ALA	5.9
43	DW	2	HIS	5.8
52	BI	70	THR	5.8
19	CT	35	TYR	5.8
52	DI	97	VAL	5.8
33	DL	143	GLU	5.8
33	BL	6	LEU	5.8
33	DL	92	LEU	5.8
51	D4	37	GLN	5.8
44	DX	39	GLN	5.8
52	BI	4	VAL	5.8

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Mol	Chain	Res	Type	RSRZ
30	BH	128	HIS	5.8
51	B4	22	VAL	5.8
52	BI	6	ALA	5.7
52	BI	115	ASP	5.7
37	DP	55	HIS	5.7
52	BI	2	LYS	5.7
36	DO	57	ALA	5.7
27	BE	145	ASP	5.7
25	DC	3	VAL	5.7
27	DE	23	PHE	5.7
52	BI	47	SER	5.7
30	BH	122	LEU	5.7
27	DE	168	ASP	5.7
12	CM	46	GLU	5.7
26	BD	186	LEU	5.7
37	BP	23	ASP	5.7
40	BS	40	ASN	5.7
52	DI	6	ALA	5.7
43	DW	79	ILE	5.7
36	DO	64	TYR	5.6
34	BM	2	LEU	5.6
50	B3	62	PRO	5.6
36	BO	56	LYS	5.6
38	DQ	96	ASP	5.6
51	D4	23	ILE	5.6
20	CB	220	VAL	5.6
42	DU	69	VAL	5.6
18	AS	14	LEU	5.6
12	AM	83	GLY	5.6
18	AS	29	PRO	5.6
29	BG	51	PHE	5.6
36	DO	61	GLN	5.6
30	DH	19	VAL	5.5
34	BM	92	TRP	5.5
39	DR	42	ALA	5.5
18	CS	2	ARG	5.5
33	BL	10	GLU	5.5
48	B1	2	LYS	5.5
26	DD	55	LYS	5.5
25	BC	238	ASN	5.5
25	BC	267	VAL	5.5
27	DE	111	GLU	5.5

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Mol	Chain	Res	Type	RSRZ
29	BG	42	VAL	5.5
8	CI	31	GLN	5.5
27	BE	5	LEU	5.5
51	B4	28	SER	5.5
18	AS	2	ARG	5.5
47	B0	56	LYS	5.5
50	B3	23	HIS	5.4
35	BN	123	GLU	5.4
27	BE	4	VAL	5.4
30	BH	149	GLU	5.4
52	BI	97	VAL	5.4
12	AM	55	LEU	5.4
52	DI	84	GLY	5.4
18	CS	78	THR	5.4
43	BW	68	PHE	5.4
43	BW	53	GLY	5.4
52	DI	132	ALA	5.4
44	DX	19	LEU	5.4
43	DW	4	LYS	5.4
19	CT	86	ALA	5.4
38	DQ	117	ALA	5.4
51	B4	25	VAL	5.4
43	DW	44	PHE	5.4
44	BX	2	LYS	5.4
18	AS	47	THR	5.4
16	AQ	27	PHE	5.4
26	DD	49	GLN	5.4
42	BU	91	LYS	5.3
36	DO	24	THR	5.3
23	DB	2402	U	5.3
33	BL	34	GLY	5.3
33	BL	4	ASN	5.3
51	D4	25	VAL	5.3
44	DX	24	GLU	5.3
19	AT	27	MET	5.3
3	AD	3	TYR	5.3
40	BS	1	MET	5.3
52	DI	53	PRO	5.3
35	DN	125	ALA	5.3
19	CT	67	HIS	5.3
5	AF	52	ASN	5.3
43	DW	3	LYS	5.3

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Mol	Chain	Res	Type	RSRZ
25	DC	209	ALA	5.2
27	DE	165	HIS	5.2
31	BJ	10	THR	5.2
26	BD	9	VAL	5.2
52	DI	52	LEU	5.2
35	BN	119	SER	5.2
21	AU	44	ARG	5.2
45	DY	58	GLU	5.2
27	BE	115	GLN	5.2
42	DU	68	ASN	5.2
36	BO	19	GLN	5.2
20	CB	123	GLY	5.2
21	CU	41	THR	5.2
43	DW	45	HIS	5.2
26	BD	133	THR	5.2
49	B2	13	ASN	5.1
33	BL	37	GLY	5.1
52	BI	3	LYS	5.1
8	AI	127	SER	5.1
6	AG	145	GLU	5.1
20	CB	124	THR	5.1
30	BH	133	GLN	5.1
47	D0	34	GLY	5.1
26	BD	64	GLU	5.1
52	DI	30	GLN	5.1
33	DL	3	LEU	5.1
36	BO	5	SER	5.1
47	D0	43	THR	5.1
27	DE	112	LEU	5.1
52	DI	111	THR	5.1
13	AN	46	LYS	5.0
52	BI	52	LEU	5.0
44	BX	27	ASN	5.0
3	AD	146	GLU	5.0
13	AN	30	ILE	5.0
26	DD	92	VAL	5.0
37	BP	49	ILE	5.0
44	DX	47	ARG	5.0
34	DM	91	TYR	5.0
12	AM	79	LEU	5.0
23	BB	2799	A	5.0
41	BT	6	ARG	5.0

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Mol	Chain	Res	Type	RSRZ
23	BB	2309	A	5.0
52	DI	136	GLY	5.0
34	BM	3	GLN	5.0
18	AS	60	PHE	5.0
50	D3	62	PRO	4.9
2	AC	98	ALA	4.9
52	BI	20	SER	4.9
30	BH	146	VAL	4.9
43	DW	67	LYS	4.9
39	BR	11	GLN	4.9
41	BT	72	GLN	4.9
41	DT	74	ILE	4.9
48	B1	53	ILE	4.9
25	BC	20	ASN	4.9
27	DE	24	ASN	4.9
30	BH	116	ARG	4.9
52	DI	85	ILE	4.9
52	BI	53	PRO	4.9
52	DI	21	PRO	4.9
25	BC	258	SER	4.9
19	CT	51	ASN	4.9
30	BH	20	ASN	4.9
39	DR	12	HIS	4.9
30	BH	127	GLU	4.9
45	BY	57	GLU	4.9
43	BW	81	ILE	4.9
18	AS	18	VAL	4.8
33	BL	12	SER	4.8
33	BL	117	THR	4.8
15	AP	82	ALA	4.8
29	DG	42	VAL	4.8
43	DW	37	VAL	4.8
52	BI	21	PRO	4.8
8	CI	128	LYS	4.8
35	BN	71	ARG	4.8
39	BR	52	PRO	4.8
33	BL	76	GLU	4.8
45	DY	57	GLU	4.8
51	B4	35	GLN	4.8
51	B4	24	ARG	4.8
25	DC	247	TRP	4.8
13	AN	20	PHE	4.8

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Mol	Chain	Res	Type	RSRZ
22	BA	88	C	4.8
41	DT	98	GLY	4.8
51	D4	38	GLY	4.8
30	BH	84	ALA	4.8
18	AS	48	ILE	4.8
41	DT	69	ARG	4.8
25	BC	4	LYS	4.8
36	DO	23	ALA	4.8
1	AA	121	U	4.8
33	BL	78	ARG	4.7
30	DH	149	GLU	4.7
36	DO	89	ASP	4.7
37	BP	20	ARG	4.7
34	BM	103	TYR	4.7
16	AQ	6	THR	4.7
30	BH	132	PHE	4.7
36	BO	89	ASP	4.7
29	BG	93	TYR	4.7
30	BH	90	LEU	4.7
16	AQ	45	VAL	4.7
27	BE	128	ALA	4.7
23	BB	654	A	4.7
36	DO	35	ILE	4.7
31	DJ	43	GLU	4.7
52	DI	25	PRO	4.7
52	DI	5	GLN	4.7
28	BF	31	GLU	4.7
41	DT	5	GLU	4.7
52	BI	5	GLN	4.7
23	DB	140	C	4.7
36	BO	61	GLN	4.7
29	BG	176	LYS	4.7
35	DN	122	ALA	4.7
17	AR	36	GLY	4.7
25	DC	9	SER	4.7
2	AC	158	GLY	4.6
23	BB	2148	G	4.6
35	BN	68	ALA	4.6
18	AS	30	LEU	4.6
30	BH	123	ARG	4.6
31	BJ	98	GLU	4.6
52	DI	99	LYS	4.6

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Mol	Chain	Res	Type	RSRZ
26	BD	17	GLU	4.6
47	B0	49	ARG	4.6
30	BH	50	ARG	4.6
39	DR	33	VAL	4.6
38	DQ	72	GLY	4.6
33	BL	52	GLY	4.6
36	BO	60	GLU	4.6
41	BT	77	ARG	4.6
27	BE	129	PRO	4.6
41	DT	37	ASP	4.6
41	DT	72	GLN	4.6
30	DH	136	SER	4.6
26	BD	132	ALA	4.6
52	DI	9	LYS	4.6
52	BI	8	VAL	4.6
26	BD	25	THR	4.6
52	DI	78	LEU	4.5
20	AB	15	PHE	4.5
23	BB	1459	G	4.5
45	BY	2	LYS	4.5
37	BP	51	ASN	4.5
30	BH	96	THR	4.5
47	D0	35	GLU	4.5
21	CU	37	TYR	4.5
34	BM	55	ARG	4.5
52	BI	60	VAL	4.5
37	BP	99	LEU	4.5
42	DU	32	LYS	4.5
7	AH	74	ILE	4.5
16	CQ	6	THR	4.5
3	AD	43	ARG	4.5
36	DO	28	VAL	4.5
42	BU	87	GLU	4.5
50	D3	36	ALA	4.5
33	BL	57	LEU	4.5
50	D3	28	LEU	4.5
52	BI	54	ILE	4.5
7	AH	110	MET	4.5
26	DD	74	GLU	4.5
51	B4	9	LYS	4.5
51	B4	15	LYS	4.5
30	BH	82	SER	4.5

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Mol	Chain	Res	Type	RSRZ
51	D4	7	VAL	4.5
52	DI	7	TYR	4.5
43	BW	77	LYS	4.5
16	AQ	20	ILE	4.4
30	BH	76	GLU	4.4
42	DU	102	ILE	4.4
43	DW	6	GLY	4.4
44	BX	39	GLN	4.4
33	BL	40	SER	4.4
33	DL	10	GLU	4.4
52	BI	48	ILE	4.4
11	AL	123	ALA	4.4
18	AS	11	ASP	4.4
29	BG	170	THR	4.4
42	BU	93	ARG	4.4
26	BD	149	ASN	4.4
48	D1	32	LYS	4.4
26	DD	33	ARG	4.4
37	BP	88	ARG	4.4
25	BC	85	ASN	4.4
44	BX	20	ASN	4.4
31	BJ	100	VAL	4.4
30	DH	72	ILE	4.4
33	BL	39	LYS	4.4
1	AA	78	A	4.4
34	DM	56	ALA	4.4
23	BB	140	C	4.4
34	DM	136	MET	4.4
41	BT	95	PHE	4.4
28	DF	73	VAL	4.3
37	DP	3	ILE	4.3
30	BH	101	ASP	4.3
51	D4	16	ILE	4.3
29	DG	169	ARG	4.3
37	BP	97	TYR	4.3
23	DB	1459	G	4.3
40	DS	109	ASP	4.3
40	BS	93	ALA	4.3
42	BU	56	GLY	4.3
44	BX	37	LEU	4.3
52	BI	7	TYR	4.3
43	BW	42	THR	4.3

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Mol	Chain	Res	Type	RSRZ
52	DI	47	SER	4.3
25	BC	43	ASN	4.3
33	DL	49	GLY	4.3
37	BP	3	ILE	4.3
47	D0	45	ASP	4.3
6	AG	38	ALA	4.3
3	AD	155	LYS	4.3
37	DP	19	PHE	4.3
27	DE	153	LEU	4.3
30	BH	104	THR	4.3
31	BJ	1	MET	4.3
52	BI	79	LEU	4.3
52	BI	116	MET	4.3
29	DG	51	PHE	4.3
31	BJ	44	TYR	4.3
39	DR	34	GLU	4.3
33	DL	53	GLY	4.3
41	DT	47	VAL	4.3
52	DI	82	ALA	4.3
50	D3	32	LEU	4.3
18	AS	46	LEU	4.3
34	DM	135	VAL	4.3
2	AC	153	SER	4.2
12	AM	95	PRO	4.2
7	CH	62	LEU	4.2
27	BE	143	LEU	4.2
1	AA	1362	A	4.2
8	AI	69	GLY	4.2
27	DE	199	MET	4.2
43	BW	74	LYS	4.2
27	DE	172	ALA	4.2
27	BE	3	LEU	4.2
30	BH	141	LYS	4.2
34	DM	1	MET	4.2
35	BN	121	LYS	4.2
7	AH	102	VAL	4.2
25	DC	85	ASN	4.2
29	DG	56	GLY	4.2
41	BT	69	ARG	4.2
25	BC	165	ALA	4.2
20	AB	199	ILE	4.2
25	BC	218	THR	4.2

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Mol	Chain	Res	Type	RSRZ
6	AG	84	TYR	4.2
20	AB	32	GLY	4.2
39	DR	77	PHE	4.2
52	DI	80	LYS	4.2
52	DI	12	VAL	4.2
52	BI	85	ILE	4.2
52	DI	135	MET	4.1
16	AQ	36	PHE	4.1
50	B3	44	ARG	4.1
50	D3	33	THR	4.1
42	BU	101	THR	4.1
35	BN	6	SER	4.1
47	B0	55	ALA	4.1
27	BE	158	PHE	4.1
33	DL	2	ARG	4.1
41	DT	13	ALA	4.1
48	D1	36	LYS	4.1
23	BB	1537	G	4.1
43	DW	11	ASN	4.1
52	DI	81	LYS	4.1
30	BH	110	VAL	4.1
41	DT	27	SER	4.1
47	B0	33	SER	4.1
6	CG	104	VAL	4.1
52	DI	100	ILE	4.1
44	DX	23	ARG	4.1
52	BI	25	PRO	4.1
1	CA	85	U	4.1
39	BR	91	GLN	4.1
39	DR	11	GLN	4.1
41	BT	4	GLU	4.1
52	DI	48	ILE	4.1
14	CO	69	LEU	4.1
25	BC	247	TRP	4.1
36	BO	92	PHE	4.1
50	B3	13	PHE	4.1
15	AP	43	ALA	4.1
52	BI	31	GLY	4.1
52	BI	11	GLN	4.1
7	AH	13	ILE	4.1
18	CS	10	ILE	4.1
44	BX	3	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
29	DG	176	LYS	4.0
34	DM	40	ARG	4.0
43	DW	19	ARG	4.0
16	AQ	7	LEU	4.0
17	CR	63	TYR	4.0
27	BE	48	THR	4.0
33	DL	58	TYR	4.0
15	AP	39	PHE	4.0
23	BB	136	G	4.0
52	DI	26	ALA	4.0
37	BP	87	ARG	4.0
33	BL	2	ARG	4.0
35	BN	20	MET	4.0
50	D3	7	ARG	4.0
17	AR	38	ILE	4.0
29	DG	172	GLU	4.0
50	D3	35	LYS	4.0
3	AD	36	ALA	4.0
52	BI	14	ALA	4.0
18	AS	36	ARG	4.0
29	BG	35	THR	4.0
37	BP	59	THR	4.0
44	BX	62	GLY	4.0
39	DR	98	ILE	4.0
38	DQ	46	TYR	4.0
30	DH	18	GLN	4.0
3	AD	154	VAL	4.0
35	DN	57	THR	4.0
47	B0	45	ASP	4.0
18	CS	65	MET	4.0
13	CN	30	ILE	4.0
41	BT	79	ASP	4.0
23	BB	2147	A	4.0
49	B2	18	PHE	4.0
26	BD	148	GLN	3.9
43	BW	3	LYS	3.9
28	DF	112	ASP	3.9
52	DI	89	SER	3.9
30	DH	12	LEU	3.9
28	BF	45	ASP	3.9
26	BD	34	VAL	3.9
1	AA	466	A	3.9

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Mol	Chain	Res	Type	RSRZ
23	DB	2799	A	3.9
27	BE	15	SER	3.9
30	BH	131	SER	3.9
8	AI	29	ILE	3.9
39	BR	92	TRP	3.9
44	DX	2	LYS	3.9
52	DI	58	ILE	3.9
52	DI	75	ALA	3.9
13	AN	40	ARG	3.9
3	AD	173	ASP	3.9
31	BJ	130	HIS	3.9
29	DG	170	THR	3.9
30	BH	18	GLN	3.9
43	BW	67	LYS	3.9
41	DT	31	VAL	3.9
1	CA	1362	A	3.9
30	BH	19	VAL	3.9
52	BI	78	LEU	3.9
42	BU	32	LYS	3.9
13	AN	27	LYS	3.9
14	CO	45	HIS	3.9
41	BT	5	GLU	3.9
20	AB	128	LEU	3.9
25	BC	10	PRO	3.9
33	DL	8	PRO	3.9
48	D1	52	LYS	3.8
50	D3	60	CYS	3.8
52	BI	86	LYS	3.8
18	CS	64	GLU	3.8
20	AB	157	PRO	3.8
34	BM	109	PRO	3.8
52	BI	87	SER	3.8
6	CG	38	ALA	3.8
52	DI	57	VAL	3.8
35	BN	72	ASP	3.8
46	DZ	2	LYS	3.8
20	AB	211	LEU	3.8
12	AM	85	TYR	3.8
41	DT	94	ASP	3.8
27	DE	144	GLU	3.8
28	DF	72	SER	3.8
37	DP	80	VAL	3.8

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Mol	Chain	Res	Type	RSRZ
41	BT	32	LEU	3.8
41	DT	96	VAL	3.8
23	BB	645	C	3.8
13	CN	34	ASN	3.8
40	BS	103	ILE	3.8
14	AO	58	MET	3.8
45	BY	1	ALA	3.8
50	B3	26	ALA	3.8
23	BB	2153	C	3.8
39	BR	26	ASP	3.8
23	BB	546	U	3.8
26	BD	147	GLY	3.8
29	DG	39	ALA	3.8
38	DQ	58	GLN	3.8
52	BI	68	PHE	3.8
30	DH	25	TYR	3.8
3	AD	28	ASP	3.8
31	DJ	1	MET	3.8
25	DC	210	ALA	3.8
17	AR	73	HIS	3.8
48	D1	31	GLU	3.8
26	DD	79	LEU	3.8
38	BQ	90	ASP	3.8
25	DC	218	THR	3.8
20	CB	38	HIS	3.8
37	BP	4	ILE	3.8
34	BM	65	ILE	3.8
24	BV	91	PHE	3.8
52	DI	37	PHE	3.8
16	AQ	9	GLY	3.8
30	DH	112	LYS	3.8
41	BT	38	ALA	3.8
12	CM	16	ILE	3.8
33	BL	19	LEU	3.8
33	BL	53	GLY	3.7
33	DL	140	GLY	3.7
41	DT	33	LYS	3.7
52	BI	95	ASP	3.7
25	DC	257	ARG	3.7
33	BL	62	PRO	3.7
20	AB	31	PHE	3.7
31	DJ	50	THR	3.7

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Mol	Chain	Res	Type	RSRZ
13	AN	21	ALA	3.7
30	BH	98	ASP	3.7
39	BR	12	HIS	3.7
33	DL	48	ARG	3.7
52	DI	70	THR	3.7
28	BF	42	ALA	3.7
13	AN	59	GLN	3.7
29	BG	21	GLN	3.7
1	CA	121	U	3.7
52	BI	108	ILE	3.7
26	BD	126	ASN	3.7
30	DH	50	ARG	3.7
37	BP	74	GLN	3.7
52	BI	128	ILE	3.7
43	DW	25	PHE	3.7
52	BI	18	ASN	3.7
42	DU	91	LYS	3.7
20	CB	204	ASP	3.7
30	BH	97	ARG	3.7
39	DR	5	PHE	3.7
28	DF	68	LYS	3.7
30	BH	87	GLU	3.7
48	B1	54	LYS	3.7
36	BO	8	ILE	3.7
37	DP	75	THR	3.7
42	BU	71	ILE	3.7
11	CL	24	GLU	3.7
52	BI	76	ALA	3.7
1	AA	79	G	3.7
39	BR	41	ILE	3.7
41	DT	91	GLN	3.7
16	AQ	55	GLY	3.7
31	DJ	49	ASP	3.7
38	DQ	91	ARG	3.7
43	BW	35	ILE	3.7
42	BU	90	LYS	3.7
19	AT	41	GLY	3.7
35	BN	7	GLY	3.7
18	AS	31	ARG	3.7
25	DC	242	HIS	3.7
6	AG	7	GLY	3.7
21	CU	35	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
30	BH	126	GLY	3.7
34	BM	125	PRO	3.7
26	DD	64	GLU	3.7
21	CU	42	THR	3.7
13	AN	5	MET	3.7
30	DH	21	VAL	3.7
6	AG	4	ARG	3.7
33	DL	59	ARG	3.7
51	D4	34	LYS	3.6
30	DH	38	PRO	3.6
35	BN	14	SER	3.6
36	BO	52	SER	3.6
37	BP	77	SER	3.6
42	BU	14	THR	3.6
18	CS	40	PHE	3.6
19	AT	34	VAL	3.6
20	CB	34	ARG	3.6
28	DF	24	VAL	3.6
41	DT	53	VAL	3.6
41	BT	25	GLU	3.6
13	AN	45	LEU	3.6
35	DN	1	MET	3.6
18	CS	48	ILE	3.6
18	CS	73	PHE	3.6
39	DR	55	ASP	3.6
13	AN	48	GLN	3.6
35	BN	1	MET	3.6
50	B3	12	ARG	3.6
39	DR	27	ILE	3.6
20	AB	212	TYR	3.6
26	BD	183	GLU	3.6
27	DE	25	GLU	3.6
30	DH	17	ASP	3.6
26	BD	150	GLN	3.6
29	DG	44	HIS	3.6
27	DE	48	THR	3.6
3	CD	28	ASP	3.6
18	CS	68	HIS	3.6
39	DR	90	ARG	3.6
41	DT	80	TRP	3.6
43	DW	82	GLU	3.6
52	DI	49	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
48	D1	48	TYR	3.6
13	AN	50	LEU	3.6
28	BF	152	ASP	3.6
41	DT	89	GLU	3.6
14	AO	16	ARG	3.6
36	DO	21	LEU	3.6
52	BI	118	GLY	3.6
52	BI	137	LEU	3.6
42	BU	45	GLN	3.6
11	CL	123	ALA	3.6
12	CM	60	ALA	3.6
42	BU	47	PRO	3.6
51	D4	32	LYS	3.6
15	AP	44	SER	3.6
1	AA	461	A	3.6
33	DL	35	HIS	3.6
50	D3	48	MET	3.6
37	DP	61	ARG	3.6
52	BI	119	ALA	3.6
18	CS	59	VAL	3.6
19	CT	38	ILE	3.6
25	BC	3	VAL	3.6
36	DO	45	SER	3.6
39	BR	71	LYS	3.6
1	CA	1031	C	3.6
36	DO	7	ARG	3.6
30	BH	109	GLU	3.6
37	BP	28	LYS	3.6
43	BW	37	VAL	3.6
52	DI	38	CYS	3.6
34	DM	18	ARG	3.5
5	CF	96	VAL	3.5
34	BM	126	ILE	3.5
13	CN	98	ALA	3.5
29	BG	150	TYR	3.5
52	DI	101	SER	3.5
41	DT	54	GLU	3.5
25	BC	224	MET	3.5
1	AA	83	C	3.5
8	CI	39	GLY	3.5
25	BC	269	ARG	3.5
2	AC	151	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
12	CM	59	VAL	3.5
28	BF	116	LEU	3.5
37	BP	61	ARG	3.5
23	BB	2104	C	3.5
46	DZ	41	HIS	3.5
27	DE	142	ALA	3.5
28	BF	131	VAL	3.5
37	BP	27	VAL	3.5
52	BI	125	THR	3.5
15	CP	47	GLU	3.5
33	BL	35	HIS	3.5
42	BU	68	ASN	3.5
28	BF	112	ASP	3.5
18	CS	60	PHE	3.5
27	BE	41	GLN	3.5
36	DO	107	ALA	3.5
23	BB	2146	C	3.5
25	BC	241	LYS	3.5
52	BI	49	GLU	3.5
27	BE	113	VAL	3.5
20	AB	68	PHE	3.5
30	DH	116	ARG	3.5
37	BP	18	SER	3.5
52	DI	79	LEU	3.5
36	BO	43	ASN	3.5
3	AD	108	ALA	3.5
30	DH	139	PHE	3.5
18	CS	1	PRO	3.5
26	BD	178	VAL	3.5
52	DI	4	VAL	3.5
52	BI	10	LEU	3.5
13	AN	68	ARG	3.5
29	DG	2	ARG	3.5
31	BJ	120	ARG	3.5
25	BC	127	ASN	3.4
30	DH	130	VAL	3.4
36	BO	3	LYS	3.4
18	CS	74	ALA	3.4
43	BW	84	GLU	3.4
13	CN	42	ASN	3.4
30	BH	83	LYS	3.4
45	BY	41	PRO	3.4

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Mol	Chain	Res	Type	RSRZ
13	AN	55	SER	3.4
34	DM	132	THR	3.4
52	DI	125	THR	3.4
27	BE	200	LEU	3.4
23	BB	2154	A	3.4
23	DB	2134	A	3.4
25	DC	269	ARG	3.4
5	CF	94	HIS	3.4
30	BH	105	ALA	3.4
16	AQ	8	GLN	3.4
15	CP	52	LEU	3.4
18	CS	70	LEU	3.4
23	BB	1538	G	3.4
30	BH	144	VAL	3.4
47	B0	41	HIS	3.4
36	BO	87	ILE	3.4
46	DZ	64	PHE	3.4
43	BW	8	SER	3.4
47	B0	34	GLY	3.4
31	BJ	99	ARG	3.4
48	D1	12	SER	3.4
1	AA	87	C	3.4
6	AG	8	GLN	3.4
6	CG	152	HIS	3.4
13	CN	46	LYS	3.4
16	CQ	3	LYS	3.4
39	BR	37	GLU	3.4
52	DI	134	SER	3.4
3	AD	175	GLY	3.4
36	BO	76	LYS	3.4
46	BZ	63	ARG	3.4
2	CC	166	TRP	3.4
26	DD	80	TRP	3.4
46	DZ	20	ASN	3.4
16	AQ	29	LYS	3.4
14	AO	55	LEU	3.4
22	BA	52	A	3.4
52	BI	139	VAL	3.4
17	CR	58	ILE	3.4
18	AS	9	PHE	3.4
20	CB	126	ASP	3.4
28	BF	129	MET	3.4

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Mol	Chain	Res	Type	RSRZ
9	AJ	34	ALA	3.4
18	AS	59	VAL	3.4
27	BE	117	ARG	3.4
36	DO	37	ALA	3.4
51	B4	7	VAL	3.4
21	AU	41	THR	3.4
41	BT	27	SER	3.4
49	D2	43	THR	3.4
52	BI	30	GLN	3.4
6	AG	79	VAL	3.4
45	BY	49	ALA	3.4
43	BW	9	THR	3.4
1	CA	1534	A	3.4
16	AQ	82	VAL	3.4
29	DG	59	ASP	3.4
33	DL	45	GLY	3.4
45	BY	38	GLU	3.4
21	CU	34	ARG	3.4
34	BM	5	LYS	3.4
37	BP	92	ARG	3.4
47	D0	25	THR	3.4
48	D1	23	THR	3.4
48	D1	43	ARG	3.4
49	B2	46	LYS	3.4
7	AH	33	VAL	3.3
41	DT	67	VAL	3.3
8	CI	64	ILE	3.3
12	CM	52	ILE	3.3
3	AD	35	GLN	3.3
12	CM	109	LYS	3.3
35	BN	15	SER	3.3
26	BD	2	ILE	3.3
2	CC	154	GLY	3.3
36	DO	19	GLN	3.3
18	AS	43	MET	3.3
26	BD	140	HIS	3.3
42	BU	69	VAL	3.3
13	CN	72	PHE	3.3
26	DD	127	PHE	3.3
44	BX	26	PHE	3.3
13	AN	23	ARG	3.3
25	DC	28	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
41	DT	68	LYS	3.3
7	AH	58	LEU	3.3
13	CN	73	LEU	3.3
41	DT	28	ASN	3.3
41	DT	50	LEU	3.3
41	DT	34	VAL	3.3
13	CN	22	LYS	3.3
19	AT	30	PHE	3.3
25	DC	156	SER	3.3
29	DG	1	SER	3.3
20	AB	34	ARG	3.3
28	BF	153	ILE	3.3
7	CH	120	LEU	3.3
19	CT	65	LEU	3.3
36	BO	106	LEU	3.3
43	DW	59	PHE	3.3
29	DG	41	GLU	3.3
37	BP	75	THR	3.3
27	BE	60	TRP	3.3
29	BG	147	LEU	3.3
34	DM	106	ASP	3.3
23	BB	138	U	3.3
25	BC	239	PHE	3.3
27	DE	141	MET	3.3
52	DI	138	VAL	3.3
20	CB	163	ILE	3.3
33	DL	93	ASN	3.3
28	DF	175	PRO	3.3
30	BH	75	LEU	3.3
25	DC	167	ASP	3.3
41	DT	10	VAL	3.3
52	DI	110	GLN	3.3
46	BZ	22	MET	3.3
15	CP	43	ALA	3.3
18	AS	13	HIS	3.3
21	CU	8	ASN	3.3
42	DU	2	ALA	3.3
8	CI	127	SER	3.3
34	DM	128	THR	3.3
28	BF	73	VAL	3.3
50	D3	53	ASP	3.3
14	AO	72	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
34	BM	83	GLY	3.3
40	DS	108	SER	3.3
41	DT	26	LYS	3.3
39	BR	25	LEU	3.3
3	AD	191	SER	3.3
37	DP	24	THR	3.3
40	BS	34	ASP	3.3
3	AD	4	LEU	3.3
12	AM	82	LEU	3.3
30	BH	121	VAL	3.3
18	AS	58	PRO	3.3
31	BJ	119	PHE	3.3
25	DC	258	SER	3.3
7	AH	125	ILE	3.3
37	BP	50	ARG	3.3
35	BN	66	ALA	3.3
18	CS	79	TYR	3.3
26	DD	68	PHE	3.3
43	DW	74	LYS	3.3
6	CG	105	GLU	3.2
8	AI	56	MET	3.2
25	DC	59	GLN	3.2
4	CE	158	LYS	3.2
3	AD	197	HIS	3.2
32	BK	122	VAL	3.2
33	BL	50	PHE	3.2
37	BP	73	PHE	3.2
42	BU	84	PHE	3.2
31	BJ	131	ASN	3.2
33	DL	28	GLY	3.2
43	BW	51	GLY	3.2
52	DI	128	ILE	3.2
9	CJ	102	LEU	3.2
12	AM	88	LEU	3.2
13	CN	18	LYS	3.2
12	AM	78	ARG	3.2
33	DL	50	PHE	3.2
37	BP	60	VAL	3.2
33	BL	144	GLU	3.2
52	BI	73	PRO	3.2
11	AL	122	LYS	3.2
27	BE	153	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
33	BL	84	LYS	3.2
26	DD	100	LEU	3.2
29	BG	148	ARG	3.2
34	BM	10	ARG	3.2
33	DL	67	THR	3.2
27	BE	22	ASP	3.2
33	BL	91	ASP	3.2
33	BL	142	ILE	3.2
20	AB	195	VAL	3.2
45	DY	2	LYS	3.2
6	AG	28	ILE	3.2
19	CT	53	MET	3.2
33	BL	27	LEU	3.2
41	DT	12	ARG	3.2
43	BW	5	ALA	3.2
48	D1	38	PHE	3.2
1	CA	1032	G	3.2
23	DB	141	G	3.2
27	BE	38	GLY	3.2
37	BP	55	HIS	3.2
28	BF	148	VAL	3.2
31	BJ	18	VAL	3.2
51	B4	30	GLU	3.2
52	DI	131	THR	3.2
34	BM	98	PRO	3.2
52	DI	139	VAL	3.2
34	BM	32	GLY	3.2
26	BD	164	GLN	3.2
34	BM	11	LYS	3.2
46	DZ	47	LYS	3.2
28	BF	151	LEU	3.2
33	DL	101	ILE	3.2
37	DP	37	LYS	3.2
41	BT	26	LYS	3.2
35	BN	102	PHE	3.2
37	BP	109	ILE	3.2
7	AH	2	MET	3.2
42	DU	48	VAL	3.2
27	DE	140	ASP	3.2
9	AJ	41	PRO	3.1
28	BF	83	PRO	3.1
39	DR	59	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
27	DE	125	SER	3.1
33	BL	54	GLN	3.1
35	BN	2	ARG	3.1
1	AA	85	U	3.1
26	BD	127	PHE	3.1
27	BE	127	GLU	3.1
28	BF	74	ALA	3.1
46	BZ	4	ASP	3.1
18	CS	29	PRO	3.1
33	BL	11	GLY	3.1
36	DO	8	ILE	3.1
20	CB	68	PHE	3.1
26	DD	125	TRP	3.1
25	BC	22	GLU	3.1
27	DE	86	ALA	3.1
36	BO	4	LYS	3.1
2	CC	153	SER	3.1
13	AN	31	SER	3.1
39	DR	39	LEU	3.1
50	D3	12	ARG	3.1
38	BQ	48	ASP	3.1
42	DU	49	PRO	3.1
4	AE	113	VAL	3.1
37	DP	30	TRP	3.1
41	DT	55	VAL	3.1
6	AG	77	ARG	3.1
32	DK	49	ARG	3.1
23	BB	2157	G	3.1
26	BD	53	GLY	3.1
26	BD	153	GLY	3.1
34	BM	4	PRO	3.1
45	BY	52	PHE	3.1
52	DI	96	LYS	3.1
27	BE	188	MET	3.1
50	D3	21	PHE	3.1
52	BI	65	SER	3.1
36	DO	30	ARG	3.1
52	BI	138	VAL	3.1
38	DQ	55	GLN	3.1
45	BY	23	LEU	3.1
42	BU	55	GLY	3.1
42	DU	34	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
7	CH	89	ASP	3.1
11	AL	97	VAL	3.1
17	AR	21	ASP	3.1
52	BI	35	MET	3.1
29	DG	160	GLY	3.1
36	DO	22	GLY	3.1
19	CT	34	VAL	3.1
13	CN	17	ASP	3.1
8	AI	16	ALA	3.1
52	DI	27	LEU	3.1
20	CB	39	ILE	3.1
45	DY	10	ARG	3.1
17	AR	39	VAL	3.1
30	BH	115	VAL	3.1
25	BC	57	HIS	3.1
38	BQ	111	LYS	3.1
27	BE	138	LEU	3.1
41	DT	57	VAL	3.1
17	AR	72	ARG	3.1
51	B4	20	ASP	3.1
8	AI	42	THR	3.1
17	AR	46	THR	3.1
36	DO	86	GLY	3.1
24	BV	59	GLU	3.1
12	CM	18	LEU	3.1
13	CN	45	LEU	3.1
26	DD	140	HIS	3.1
45	DY	29	ARG	3.1
19	CT	60	GLN	3.0
29	BG	130	ILE	3.0
46	BZ	5	ILE	3.0
31	DJ	102	GLU	3.0
36	BO	103	VAL	3.0
6	CG	29	LEU	3.0
12	CM	55	LEU	3.0
20	AB	153	MET	3.0
24	DV	56	PHE	3.0
8	AI	64	ILE	3.0
19	AT	66	ILE	3.0
37	BP	62	LYS	3.0
43	DW	18	LYS	3.0
19	AT	12	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
28	BF	30	VAL	3.0
5	AF	88	MET	3.0
20	AB	100	LEU	3.0
20	CB	56	LEU	3.0
41	DT	87	LEU	3.0
30	BH	100	ALA	3.0
41	BT	82	LYS	3.0
23	DB	2147	A	3.0
25	BC	212	TRP	3.0
27	DE	155	GLU	3.0
30	DH	76	GLU	3.0
1	AA	81	A	3.0
29	BG	161	VAL	3.0
41	DT	36	LYS	3.0
48	B1	49	LYS	3.0
50	B3	9	ALA	3.0
2	CC	165	GLU	3.0
12	AM	32	ILE	3.0
19	CT	74	HIS	3.0
35	BN	8	ARG	3.0
48	D1	27	ARG	3.0
10	AK	49	SER	3.0
25	DC	256	THR	3.0
26	BD	16	THR	3.0
23	BB	2139	U	3.0
43	DW	78	PHE	3.0
21	AU	3	ILE	3.0
26	DD	59	ARG	3.0
29	BG	92	GLY	3.0
50	D3	19	GLY	3.0
15	AP	71	VAL	3.0
24	BV	34	LYS	3.0
39	DR	87	GLN	3.0
28	BF	161	SER	3.0
31	BJ	89	PHE	3.0
8	CI	8	THR	3.0
31	BJ	3	THR	3.0
31	DJ	10	THR	3.0
35	BN	36	THR	3.0
29	BG	120	ILE	3.0
10	AK	128	VAL	3.0
18	CS	50	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
31	BJ	47	HIS	3.0
33	DL	62	PRO	3.0
35	BN	48	VAL	3.0
51	D4	18	LYS	3.0
37	DP	40	GLN	3.0
41	BT	7	LEU	3.0
16	AQ	14	ASP	3.0
48	D1	39	ASP	3.0
25	DC	20	ASN	3.0
3	AD	42	ALA	3.0
37	BP	98	TYR	3.0
3	AD	6	PRO	3.0
31	BJ	40	HIS	3.0
52	DI	104	GLN	3.0
33	BL	126	ARG	3.0
40	DS	1	MET	3.0
49	B2	35	ARG	3.0
17	AR	20	ILE	3.0
19	CT	62	ALA	3.0
20	CB	35	ASN	3.0
48	D1	8	ILE	3.0
34	BM	33	LEU	3.0
39	DR	92	TRP	3.0
52	BI	66	PHE	3.0
6	AG	6	ILE	3.0
27	BE	155	GLU	3.0
27	DE	22	ASP	3.0
41	BT	10	VAL	3.0
28	BF	35	LEU	3.0
29	BG	155	PRO	3.0
33	DL	115	GLU	3.0
27	BE	49	ARG	3.0
38	BQ	50	ARG	3.0
28	DF	174	PHE	3.0
48	D1	16	THR	3.0
52	BI	112	LYS	3.0
18	CS	43	MET	3.0
6	CG	66	GLU	3.0
25	BC	209	ALA	3.0
29	DG	89	VAL	3.0
27	BE	72	SER	3.0
52	BI	134	SER	3.0

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Mol	Chain	Res	Type	RSRZ
13	AN	9	GLU	3.0
37	BP	22	GLY	2.9
44	BX	5	GLU	3.0
30	BH	111	ALA	2.9
29	DG	55	ASP	2.9
33	BL	66	PHE	2.9
4	CE	115	GLU	2.9
22	DA	52	A	2.9
31	DJ	80	HIS	2.9
47	B0	51	ARG	2.9
52	DI	116	MET	2.9
7	CH	129	ALA	2.9
40	BS	66	ILE	2.9
9	AJ	42	LEU	2.9
12	CM	82	LEU	2.9
15	AP	80	LYS	2.9
31	BJ	9	GLU	2.9
35	DN	121	LYS	2.9
36	DO	3	LYS	2.9
1	CA	461	A	2.9
28	BF	142	TYR	2.9
51	D4	3	VAL	2.9
36	DO	97	PHE	2.9
42	BU	86	PHE	2.9
14	AO	61	GLN	2.9
30	BH	73	ASN	2.9
33	DL	68	SER	2.9
41	DT	44	LYS	2.9
7	AH	128	VAL	2.9
13	AN	51	PRO	2.9
46	BZ	27	THR	2.9
3	AD	145	ARG	2.9
23	BB	613	A	2.9
29	BG	159	LYS	2.9
42	DU	88	ASP	2.9
23	DB	62	U	2.9
30	BH	74	ALA	2.9
30	DH	28	ASN	2.9
47	B0	53	VAL	2.9
19	AT	35	TYR	2.9
25	BC	257	ARG	2.9
25	DC	4	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
31	DJ	2	LYS	2.9
52	DI	74	PRO	2.9
26	BD	14	ILE	2.9
33	BL	55	MET	2.9
52	DI	95	ASP	2.9
42	BU	44	HIS	2.9
43	BW	43	LYS	2.9
52	BI	26	ALA	2.9
52	DI	76	ALA	2.9
25	BC	156	SER	2.9
48	B1	12	SER	2.9
16	CQ	26	ARG	2.9
41	DT	8	LEU	2.9
8	CI	15	ALA	2.9
28	DF	55	ASP	2.9
23	BB	144	A	2.9
26	DD	126	ASN	2.9
33	BL	25	SER	2.9
19	CT	78	LEU	2.9
29	BG	91	VAL	2.9
2	AC	99	GLN	2.9
5	CF	80	PHE	2.9
30	BH	102	ALA	2.9
35	BN	19	ALA	2.9
39	DR	45	GLU	2.9
46	DZ	18	CYS	2.9
41	BT	96	VAL	2.9
45	BY	24	LEU	2.9
8	AI	38	PHE	2.9
34	DM	13	HIS	2.9
36	DO	16	ARG	2.9
8	AI	62	LEU	2.9
27	DE	110	SER	2.9
50	D3	46	LYS	2.9
30	DH	3	VAL	2.9
34	BM	124	LEU	2.9
39	BR	19	THR	2.9
29	DG	58	ALA	2.9
26	DD	17	GLU	2.9
33	DL	51	GLU	2.9
35	BN	112	TYR	2.9
35	DN	45	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
12	CM	43	LYS	2.9
8	AI	124	PRO	2.9
27	DE	55	SER	2.8
52	BI	69	VAL	2.8
31	BJ	88	THR	2.8
1	CA	1042	A	2.8
41	DT	32	LEU	2.8
52	BI	84	GLY	2.8
42	BU	58	VAL	2.8
48	D1	18	HIS	2.8
24	BV	1	MET	2.8
6	CG	8	GLN	2.8
11	AL	25	ALA	2.8
52	BI	27	LEU	2.8
45	DY	37	ARG	2.8
46	BZ	67	PRO	2.8
12	CM	31	ALA	2.8
49	D2	45	SER	2.8
8	CI	65	THR	2.8
26	BD	52	THR	2.8
46	DZ	13	THR	2.8
52	BI	72	THR	2.8
2	AC	82	ASP	2.8
8	CI	40	ARG	2.8
25	BC	202	ARG	2.8
40	DS	66	ILE	2.8
52	DI	8	VAL	2.8
7	CH	44	PHE	2.8
23	DB	1205	A	2.8
33	DL	42	SER	2.8
34	BM	64	TRP	2.8
37	DP	74	GLN	2.8
28	DF	33	ILE	2.8
37	DP	60	VAL	2.8
18	CS	58	PRO	2.8
29	BG	11	PRO	2.8
29	DG	53	PRO	2.8
30	DH	128	HIS	2.8
26	DD	75	ALA	2.8
26	BD	130	GLN	2.8
33	DL	102	GLY	2.8
37	BP	114	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
39	DR	80	ARG	2.8
46	BZ	33	ASN	2.8
24	BV	63	ILE	2.8
7	CH	9	MET	2.8
6	CG	20	GLU	2.8
26	BD	28	GLU	2.8
50	B3	42	HIS	2.8
3	AD	153	ARG	2.8
19	AT	9	ARG	2.8
20	AB	159	ALA	2.8
30	DH	140	ALA	2.8
33	DL	70	LYS	2.8
38	DQ	51	GLN	2.8
12	AM	29	SER	2.8
33	BL	103	ILE	2.8
52	BI	100	ILE	2.8
46	DZ	45	THR	2.8
2	AC	159	ALA	2.8
3	AD	20	LEU	2.8
1	AA	82	G	2.8
29	BG	160	GLY	2.8
43	DW	39	GLN	2.8
5	AF	51	ILE	2.8
7	CH	78	SER	2.8
9	AJ	36	VAL	2.8
37	DP	63	ILE	2.8
45	BY	13	ILE	2.8
50	D3	13	PHE	2.8
37	DP	65	ASN	2.8
9	CJ	34	ALA	2.8
15	AP	6	LEU	2.8
28	DF	15	LEU	2.8
46	DZ	9	TYR	2.8
26	BD	170	VAL	2.8
27	BE	186	VAL	2.8
29	BG	105	SER	2.8
42	DU	99	SER	2.8
23	BB	1074	G	2.8
50	B3	24	LYS	2.8
51	D4	13	ASN	2.8
39	DR	29	THR	2.8
16	AQ	43	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
48	B1	18	HIS	2.8
52	DI	10	LEU	2.8
8	CI	18	VAL	2.8
26	BD	146	ILE	2.8
48	B1	6	GLU	2.8
30	DH	43	ASN	2.8
25	DC	10	PRO	2.8
39	DR	83	TYR	2.8
2	AC	195	ILE	2.8
3	AD	120	LYS	2.8
7	AH	100	ILE	2.8
52	DI	11	GLN	2.8
52	DI	54	ILE	2.8
23	DB	846	U	2.8
39	DR	62	GLU	2.8
18	AS	65	MET	2.8
42	BU	40	LEU	2.8
39	BR	56	GLY	2.8
41	DT	97	GLY	2.8
9	CJ	51	VAL	2.8
7	AH	17	GLN	2.7
13	AN	70	HIS	2.8
19	CT	63	LYS	2.8
33	DL	85	VAL	2.8
39	DR	79	ARG	2.8
42	DU	76	THR	2.8
46	DZ	59	ARG	2.8
48	D1	46	VAL	2.8
2	AC	179	ALA	2.7
13	AN	57	SER	2.7
20	CB	147	LEU	2.7
41	DT	61	LEU	2.7
33	DL	141	LYS	2.7
35	DN	59	SER	2.7
37	BP	94	ALA	2.7
51	B4	2	LYS	2.7
36	DO	103	VAL	2.7
2	AC	9	ILE	2.7
6	AG	37	THR	2.7
18	CS	62	THR	2.7
25	DC	190	THR	2.7
31	DJ	86	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
35	BN	16	HIS	2.7
47	D0	22	THR	2.7
52	DI	29	GLN	2.7
23	BB	2402	U	2.7
50	D3	44	ARG	2.7
23	BB	2138	G	2.7
44	BX	34	SER	2.7
18	AS	42	ASN	2.7
25	BC	114	GLN	2.7
27	BE	17	THR	2.7
5	CF	86	ARG	2.7
18	AS	17	LYS	2.7
34	DM	92	TRP	2.7
42	DU	6	ARG	2.7
52	DI	35	MET	2.7
50	B3	57	VAL	2.7
19	CT	52	GLU	2.7
8	AI	128	LYS	2.7
18	AS	68	HIS	2.7
49	B2	37	LYS	2.7
52	DI	124	MET	2.7
17	AR	63	TYR	2.7
25	BC	154	ALA	2.7
29	BG	164	ALA	2.7
30	DH	67	ALA	2.7
31	BJ	75	TYR	2.7
15	CP	10	GLY	2.7
20	CB	69	VAL	2.7
28	DF	30	VAL	2.7
52	DI	43	ALA	2.7
27	BE	123	LYS	2.7
36	DO	40	ILE	2.7
7	AH	5	PRO	2.7
30	BH	117	LEU	2.7
32	DK	14	SER	2.7
45	DY	11	SER	2.7
46	DZ	22	MET	2.7
3	AD	107	GLY	2.7
26	BD	131	ASP	2.7
33	DL	107	PHE	2.7
36	DO	92	PHE	2.7
41	BT	57	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
18	AS	10	ILE	2.7
27	BE	40	ARG	2.7
7	AH	28	SER	2.7
25	DC	246	PRO	2.7
33	DL	56	PRO	2.7
52	DI	22	PRO	2.7
23	BB	1728	C	2.7
28	DF	75	GLY	2.7
41	BT	65	GLY	2.7
52	BI	91	LYS	2.7
3	AD	1	ALA	2.7
52	DI	56	VAL	2.7
20	AB	66	ILE	2.7
20	AB	161	PHE	2.7
29	BG	157	LYS	2.7
52	BI	55	PRO	2.7
49	B2	28	ARG	2.7
52	DI	71	LYS	2.7
2	AC	75	VAL	2.7
36	BO	78	VAL	2.7
42	BU	52	ASN	2.7
29	BG	59	ASP	2.7
49	B2	43	THR	2.7
25	BC	116	GLN	2.7
34	BM	22	GLN	2.7
37	DP	92	ARG	2.7
46	DZ	48	GLN	2.7
13	AN	71	GLY	2.7
49	D2	16	HIS	2.7
50	B3	30	HIS	2.7
50	D3	20	GLY	2.7
52	DI	55	PRO	2.7
20	AB	95	TRP	2.7
27	DE	122	GLU	2.7
43	BW	80	SER	2.7
48	D1	13	SER	2.7
1	CA	1033	G	2.7
25	DC	139	THR	2.7
10	CK	125	LYS	2.7
14	AO	34	GLN	2.7
25	BC	51	ARG	2.7
34	BM	130	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
10	AK	88	PRO	2.7
15	AP	68	SER	2.7
18	CS	13	HIS	2.7
34	BM	101	VAL	2.7
24	BV	4	ILE	2.7
39	BR	42	ALA	2.7
14	AO	66	LEU	2.7
29	BG	113	ASP	2.7
29	DG	38	ASP	2.7
33	BL	3	LEU	2.7
43	BW	11	ASN	2.7
49	B2	41	ARG	2.7
44	BX	8	GLU	2.7
28	BF	7	TYR	2.6
30	BH	140	ALA	2.7
2	CC	36	PHE	2.6
12	AM	84	CYS	2.6
52	DI	67	THR	2.6
3	AD	151	GLN	2.6
4	CE	11	GLN	2.6
18	AS	55	GLN	2.6
20	CB	162	VAL	2.6
18	AS	28	LYS	2.6
26	BD	54	ALA	2.6
44	DX	42	LEU	2.6
2	AC	18	ASN	2.6
25	DC	44	ASN	2.6
2	AC	154	GLY	2.6
19	CT	70	LYS	2.6
27	BE	136	GLN	2.6
29	BG	101	VAL	2.6
33	BL	102	GLY	2.6
48	D1	41	VAL	2.6
52	BI	28	GLY	2.6
52	BI	32	VAL	2.6
18	CS	39	ILE	2.6
44	DX	21	LEU	2.6
3	AD	22	SER	2.6
30	DH	29	PHE	2.6
41	DT	95	PHE	2.6
52	DI	20	SER	2.6
7	AH	71	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
27	DE	184	ASP	2.6
41	BT	64	LYS	2.6
42	BU	60	LYS	2.6
52	DI	117	THR	2.6
12	AM	114	PRO	2.6
29	DG	106	LEU	2.6
23	DB	101	A	2.6
8	AI	92	SER	2.6
13	AN	54	SER	2.6
47	B0	28	SER	2.6
20	AB	8	MET	2.6
21	AU	35	GLU	2.6
34	DM	38	ARG	2.6
41	BT	85	VAL	2.6
34	DM	60	GLN	2.6
42	DU	7	ASP	2.6
52	DI	106	GLN	2.6
5	AF	6	ILE	2.6
14	AO	21	THR	2.6
28	BF	65	LEU	2.6
35	BN	70	THR	2.6
1	CA	1043	G	2.6
5	AF	62	MET	2.6
29	DG	3	VAL	2.6
1	AA	1317	C	2.6
20	AB	193	ASP	2.6
28	BF	56	LEU	2.6
20	AB	27	LYS	2.6
31	DJ	110	PRO	2.6
49	D2	37	LYS	2.6
11	CL	93	ARG	2.6
10	CK	128	VAL	2.6
25	DC	162	GLN	2.6
52	BI	127	SER	2.6
52	DI	108	ILE	2.6
10	CK	108	ASN	2.6
8	CI	126	PHE	2.6
9	AJ	50	THR	2.6
25	DC	160	TYR	2.6
34	BM	66	ARG	2.6
38	DQ	87	VAL	2.6
19	CT	82	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
33	BL	135	ILE	2.6
39	DR	60	LYS	2.6
41	DT	48	GLN	2.6
7	AH	115	ALA	2.6
11	CL	85	ARG	2.6
8	CI	42	THR	2.6
25	DC	199	HIS	2.6
26	DD	204	LYS	2.6
7	CH	125	ILE	2.6
17	CR	22	TYR	2.6
21	AU	37	TYR	2.6
27	BE	171	ASP	2.6
31	DJ	4	PHE	2.6
41	DT	35	ALA	2.6
43	BW	44	PHE	2.6
9	AJ	62	ARG	2.6
20	AB	156	LEU	2.6
46	DZ	6	HIS	2.6
36	BO	9	ARG	2.6
52	BI	135	MET	2.6
23	BB	62	U	2.6
8	CI	37	TYR	2.6
39	BR	93	PHE	2.6
9	AJ	91	ASP	2.6
26	BD	56	LYS	2.6
33	DL	109	LYS	2.6
20	AB	35	ASN	2.6
42	BU	41	VAL	2.6
52	BI	51	GLY	2.6
39	DR	89	HIS	2.6
34	BM	31	PHE	2.6
6	AG	144	ALA	2.6
48	D1	44	GLN	2.6
52	DI	40	ALA	2.6
3	AD	44	LYS	2.6
3	AD	196	GLU	2.6
14	AO	26	VAL	2.6
41	DT	6	ARG	2.6
43	DW	63	ASP	2.6
4	AE	35	LEU	2.6
39	BR	43	ASN	2.6
52	DI	105	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
2	CC	189	HIS	2.6
8	CI	83	THR	2.6
45	DY	8	GLN	2.5
42	BU	85	ARG	2.5
52	DI	69	VAL	2.5
9	AJ	63	ASP	2.5
17	CR	71	ASP	2.5
26	BD	18	ASP	2.5
11	CL	95	HIS	2.5
36	DO	65	THR	2.5
37	DP	110	LYS	2.5
17	AR	22	TYR	2.5
31	DJ	15	TRP	2.5
19	AT	28	ARG	2.5
1	AA	1441	A	2.5
21	CU	12	ASP	2.5
16	AQ	60	ILE	2.5
18	AS	39	ILE	2.5
46	DZ	44	PHE	2.5
17	AR	19	GLU	2.5
43	BW	57	THR	2.5
45	BY	8	GLN	2.5
18	CS	61	VAL	2.5
18	CS	46	LEU	2.5
30	DH	35	LYS	2.5
7	CH	1	SER	2.5
9	AJ	76	ILE	2.5
16	AQ	37	ILE	2.5
38	BQ	35	PHE	2.5
52	DI	16	MET	2.5
1	CA	1441	A	2.5
7	CH	123	GLU	2.5
31	BJ	80	HIS	2.5
29	BG	24	THR	2.5
42	BU	66	VAL	2.5
47	B0	25	THR	2.5
6	AG	29	LEU	2.5
28	BF	144	LYS	2.5
15	CP	38	PHE	2.5
34	BM	117	PHE	2.5
4	CE	12	GLU	2.5
18	AS	79	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	AA	1493	A	2.5
6	AG	80	GLY	2.5
26	DD	42	ASN	2.5
35	DN	25	ALA	2.5
41	BT	35	ALA	2.5
8	AI	20	ILE	2.5
8	AI	27	ILE	2.5
27	BE	199	MET	2.5
52	DI	73	PRO	2.5
6	AG	5	VAL	2.5
25	BC	59	GLN	2.5
30	DH	36	ALA	2.5
50	D3	27	ASN	2.5
28	DF	67	THR	2.5
31	DJ	88	THR	2.5
37	BP	42	PHE	2.5
52	DI	41	PHE	2.5
30	BH	60	GLU	2.5
42	DU	20	LYS	2.5
51	D4	8	LYS	2.5
6	CG	60	ALA	2.5
30	DH	82	SER	2.5
38	DQ	112	ALA	2.5
26	DD	36	GLN	2.5
8	CI	56	MET	2.5
9	AJ	27	GLU	2.5
18	AS	41	PRO	2.5
29	DG	167	VAL	2.5
38	DQ	32	ARG	2.5
40	DS	23	LEU	2.5
41	DT	73	ARG	2.5
47	D0	48	TYR	2.5
44	DX	40	SER	2.5
35	BN	3	HIS	2.5
31	BJ	4	PHE	2.5
34	BM	133	LYS	2.5
42	DU	11	ILE	2.5
48	B1	50	GLU	2.5
31	BJ	87	ALA	2.5
37	BP	85	VAL	2.5
44	DX	48	ARG	2.5
7	CH	39	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
7	CH	127	TYR	2.5
21	CU	28	LEU	2.5
27	BE	192	ALA	2.5
28	BF	141	ASP	2.5
33	BL	73	ILE	2.5
30	BH	70	GLU	2.5
18	AS	80	ARG	2.5
26	DD	197	THR	2.5
37	DP	20	ARG	2.5
8	AI	47	VAL	2.5
23	BB	2140	G	2.5
26	BD	100	LEU	2.5
26	DD	76	GLY	2.5
29	DG	161	VAL	2.5
52	BI	123	ALA	2.5
13	AN	22	LYS	2.5
11	CL	81	ILE	2.5
19	CT	31	ILE	2.5
42	BU	11	ILE	2.5
6	CG	77	ARG	2.5
28	BF	86	CYS	2.5
15	AP	19	VAL	2.5
24	BV	61	LEU	2.5
28	BF	154	THR	2.5
26	BD	116	LYS	2.5
33	BL	75	ALA	2.5
37	DP	36	LYS	2.5
9	AJ	64	GLN	2.5
33	BL	64	PHE	2.5
7	CH	74	ILE	2.5
41	DT	30	ILE	2.5
28	BF	37	MET	2.4
25	BC	225	ASN	2.4
30	DH	75	LEU	2.4
39	DR	85	LYS	2.4
42	DU	89	GLY	2.4
52	DI	33	ASN	2.4
3	AD	194	ILE	2.4
7	AH	6	ILE	2.4
26	DD	94	GLN	2.4
26	DD	128	ARG	2.4
39	BR	86	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
40	BS	74	ILE	2.4
21	AU	4	LYS	2.4
30	DH	110	VAL	2.4
37	BP	46	VAL	2.4
41	BT	34	VAL	2.4
41	BT	36	LYS	2.4
47	D0	28	SER	2.4
49	B2	8	SER	2.4
26	DD	26	VAL	2.4
37	BP	48	ALA	2.4
18	AS	75	PRO	2.4
30	DH	133	GLN	2.4
30	DH	137	GLU	2.4
29	BG	3	VAL	2.4
33	DL	47	ARG	2.4
46	BZ	9	TYR	2.4
4	CE	130	THR	2.4
5	CF	51	ILE	2.4
8	CI	20	ILE	2.4
27	BE	43	THR	2.4
28	BF	155	ILE	2.4
48	D1	21	THR	2.4
25	DC	21	PRO	2.4
33	BL	36	LYS	2.4
9	AJ	10	LEU	2.4
9	AJ	38	GLY	2.4
18	AS	70	LEU	2.4
27	DE	143	LEU	2.4
31	DJ	40	HIS	2.4
41	BT	58	VAL	2.4
36	DO	52	SER	2.4
41	BT	84	TYR	2.4
2	AC	181	ILE	2.4
20	CB	150	ILE	2.4
37	DP	47	ILE	2.4
39	BR	45	GLU	2.4
52	BI	80	LYS	2.4
38	DQ	43	GLN	2.4
8	CI	48	ARG	2.4
32	BK	78	ARG	2.4
19	AT	19	HIS	2.4
3	AD	147	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
24	BV	82	TYR	2.4
42	DU	95	PHE	2.4
2	CC	56	ILE	2.4
12	CM	32	ILE	2.4
26	DD	88	GLU	2.4
46	BZ	10	GLU	2.4
46	DZ	5	ILE	2.4
52	BI	34	ILE	2.4
6	CG	108	ARG	2.4
8	AI	39	GLY	2.4
24	DV	72	VAL	2.4
39	DR	68	ARG	2.4
40	DS	110	ARG	2.4
52	BI	67	THR	2.4
30	DH	13	GLY	2.4
2	AC	92	ASP	2.4
8	AI	19	PHE	2.4
45	BY	19	HIS	2.4
11	CL	37	TYR	2.4
2	AC	45	GLU	2.4
30	DH	113	SER	2.4
13	AN	58	ARG	2.4
14	CO	24	THR	2.4
19	CT	57	VAL	2.4
29	DG	40	VAL	2.4
39	DR	64	VAL	2.4
52	BI	45	THR	2.4
52	DI	77	VAL	2.4
13	AN	53	ASP	2.4
23	DB	1460	U	2.4
42	DU	44	HIS	2.4
15	CP	57	ILE	2.4
30	DH	99	ILE	2.4
13	AN	65	GLN	2.4
10	AK	87	GLY	2.4
3	AD	34	GLU	2.4
25	DC	57	HIS	2.4
26	DD	13	ARG	2.4
37	BP	71	ARG	2.4
51	D4	36	ARG	2.4
12	CM	38	ILE	2.4
1	AA	995	C	2.4

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Mol	Chain	Res	Type	RSRZ
4	AE	21	SER	2.4
12	AM	80	MET	2.4
28	BF	111	ARG	2.4
30	DH	73	ASN	2.4
30	BH	36	ALA	2.4
36	DO	29	HIS	2.4
8	AI	51	LEU	2.4
12	CM	12	LYS	2.4
17	AR	66	LEU	2.4
37	BP	95	LYS	2.4
38	DQ	84	LYS	2.4
3	AD	11	SER	2.4
9	AJ	74	VAL	2.4
12	AM	74	MET	2.4
1	AA	1533	C	2.4
52	BI	38	CYS	2.4
38	BQ	88	GLU	2.4
42	DU	52	ASN	2.4
48	D1	20	TYR	2.4
7	CH	124	ILE	2.4
12	AM	38	ILE	2.4
26	DD	2	ILE	2.4
51	B4	8	LYS	2.4
14	AO	20	ASP	2.4
52	BI	46	ASP	2.4
3	AD	5	GLY	2.4
20	AB	123	GLY	2.4
33	BL	24	GLY	2.4
9	AJ	68	ARG	2.4
34	BM	97	GLN	2.4
38	BQ	58	GLN	2.4
41	DT	3	ARG	2.4
8	CI	16	ALA	2.4
9	CJ	12	ALA	2.4
12	CM	4	ALA	2.4
26	DD	48	ILE	2.4
44	DX	63	ALA	2.4
8	CI	117	LEU	2.4
7	CH	103	VAL	2.3
18	CS	45	GLY	2.3
33	BL	124	GLY	2.3
42	BU	80	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
52	DI	32	VAL	2.3
37	BP	9	GLN	2.3
1	CA	84	U	2.3
27	BE	125	SER	2.3
3	AD	27	ILE	2.3
11	AL	12	ALA	2.3
36	BO	36	TYR	2.3
30	BH	6	LEU	2.3
43	BW	66	VAL	2.3
52	BI	12	VAL	2.3
27	BE	7	ASP	2.3
46	DZ	8	LYS	2.3
50	D3	22	LYS	2.3
1	AA	205	A	2.3
23	BB	1509	A	2.3
14	AO	1	SER	2.3
11	AL	11	ARG	2.3
27	BE	59	PRO	2.3
4	AE	72	ASN	2.3
25	BC	126	GLY	2.3
26	DD	32	ASN	2.3
13	AN	3	GLN	2.3
29	BG	162	ARG	2.3
31	DJ	46	PRO	2.3
35	BN	98	LEU	2.3
36	BO	21	LEU	2.3
52	BI	121	ILE	2.3
25	BC	243	PRO	2.3
8	CI	46	VAL	2.3
16	AQ	57	VAL	2.3
26	BD	166	GLY	2.3
41	BT	71	GLY	2.3
2	CC	21	TRP	2.3
16	CQ	8	GLN	2.3
28	BF	157	THR	2.3
26	BD	43	ASP	2.3
46	DZ	31	ASP	2.3
12	CM	1	ALA	2.3
23	BB	1726	C	2.3
39	BR	79	ARG	2.3
34	DM	103	TYR	2.3
45	DY	28	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
42	BU	12	VAL	2.3
9	AJ	28	THR	2.3
19	AT	17	ARG	2.3
35	BN	9	GLN	2.3
39	DR	86	GLN	2.3
41	BT	80	TRP	2.3
52	DI	59	THR	2.3
2	AC	23	ALA	2.3
3	AD	122	ILE	2.3
6	CG	119	LEU	2.3
20	AB	122	ASP	2.3
16	CQ	81	ALA	2.3
32	BK	77	ILE	2.3
41	BT	13	ALA	2.3
41	DT	11	LEU	2.3
52	DI	62	ALA	2.3
7	CH	122	GLY	2.3
42	BU	94	PHE	2.3
52	BI	136	GLY	2.3
33	BL	86	GLU	2.3
1	AA	80	A	2.3
14	AO	38	LEU	2.3
14	AO	69	LEU	2.3
30	BH	125	THR	2.3
52	DI	2	LYS	2.3
30	DH	94	ILE	2.3
52	DI	17	ALA	2.3
8	CI	38	PHE	2.3
27	DE	42	GLY	2.3
13	CN	52	ARG	2.3
48	B1	5	ARG	2.3
10	AK	79	LYS	2.3
44	DX	41	HIS	2.3
2	AC	136	ALA	2.3
26	BD	203	VAL	2.3
28	DF	137	PHE	2.3
6	CG	69	ARG	2.3
33	DL	20	GLY	2.3
33	DL	34	GLY	2.3
12	CM	9	PRO	2.3
21	AU	23	GLU	2.3
37	BP	43	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
52	BI	9	LYS	2.3
7	CH	75	GLN	2.3
18	CS	30	LEU	2.3
35	BN	51	LEU	2.3
44	BX	14	LEU	2.3
9	AJ	8	ILE	2.3
7	AH	114	ALA	2.3
3	AD	29	THR	2.3
5	CF	91	ARG	2.3
9	CJ	50	THR	2.3
26	DD	203	VAL	2.3
33	BL	21	ARG	2.3
37	BP	72	VAL	2.3
52	DI	18	ASN	2.3
28	BF	64	PRO	2.3
6	AG	46	LEU	2.3
13	AN	15	LEU	2.3
12	AM	92	ARG	2.3
20	CB	17	HIS	2.3
45	BY	31	ILE	2.3
31	DJ	6	ALA	2.3
42	DU	33	VAL	2.3
25	DC	178	GLY	2.3
6	AG	82	SER	2.3
6	CG	41	ILE	2.3
10	CK	126	ARG	2.3
20	AB	163	ILE	2.3
24	BV	56	PHE	2.3
28	BF	72	SER	2.3
33	BL	111	ILE	2.3
33	DL	18	ARG	2.3
35	BN	64	ARG	2.3
9	AJ	61	ALA	2.3
11	AL	110	LYS	2.3
27	BE	132	LYS	2.3
37	BP	19	PHE	2.3
32	BK	34	GLY	2.3
34	BM	37	GLY	2.3
37	BP	30	TRP	2.3
8	AI	30	ASN	2.3
19	CT	58	ASP	2.3
27	DE	116	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
41	DT	60	THR	2.3
4	CE	114	LEU	2.3
7	AH	9	MET	2.3
19	CT	66	ILE	2.3
3	AD	204	SER	2.3
12	AM	59	VAL	2.3
13	AN	79	SER	2.3
18	CS	22	VAL	2.3
29	BG	73	SER	2.3
4	AE	90	GLY	2.2
6	AG	81	GLY	2.2
16	CQ	51	GLU	2.2
32	BK	21	CYS	2.2
18	AS	32	THR	2.2
7	AH	27	PRO	2.2
7	AH	92	PRO	2.2
15	AP	67	ILE	2.2
50	D3	38	LYS	2.2
24	DV	91	PHE	2.2
6	CG	64	ALA	2.2
25	BC	250	GLN	2.2
34	BM	67	VAL	2.2
36	BO	28	VAL	2.2
37	BP	45	VAL	2.2
30	DH	107	GLY	2.2
36	BO	95	SER	2.2
42	DU	30	SER	2.2
8	AI	112	ARG	2.2
43	BW	40	ARG	2.2
51	B4	19	ARG	2.2
26	BD	55	LYS	2.2
8	AI	65	THR	2.2
25	BC	222	THR	2.2
31	DJ	19	ASP	2.2
44	DX	30	MET	2.2
30	DH	121	VAL	2.2
31	DJ	89	PHE	2.2
47	B0	8	THR	2.2
44	BX	11	VAL	2.2
28	DF	44	ALA	2.2
36	BO	79	ALA	2.2
43	DW	51	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
3	AD	103	ARG	2.2
25	BC	37	SER	2.2
19	AT	65	LEU	2.2
24	BV	83	LYS	2.2
37	BP	13	LYS	2.2
39	DR	25	LEU	2.2
15	CP	36	VAL	2.2
18	CS	47	THR	2.2
30	DH	47	PHE	2.2
45	BY	40	THR	2.2
46	DZ	43	PHE	2.2
52	DI	19	PRO	2.2
2	CC	70	ALA	2.2
5	CF	95	ALA	2.2
13	CN	23	ARG	2.2
14	AO	57	ARG	2.2
17	AR	60	ARG	2.2
23	BB	2141	G	2.2
23	DB	1537	G	2.2
30	DH	51	ARG	2.2
30	DH	81	ALA	2.2
52	BI	130	GLY	2.2
3	AD	10	LEU	2.2
12	CM	29	SER	2.2
18	AS	56	HIS	2.2
19	AT	18	LYS	2.2
41	DT	15	HIS	2.2
14	CO	2	LEU	2.2
27	DE	138	LEU	2.2
2	AC	150	VAL	2.2
32	BK	39	ILE	2.2
2	AC	77	GLY	2.2
18	CS	41	PRO	2.2
19	AT	23	ARG	2.2
21	CU	20	ARG	2.2
31	DJ	13	ARG	2.2
33	BL	69	ARG	2.2
31	DJ	7	LYS	2.2
34	DM	4	PRO	2.2
35	BN	69	ARG	2.2
20	AB	215	ALA	2.2
29	BG	12	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
36	DO	53	THR	2.2
36	DO	88	LYS	2.2
26	BD	27	ILE	2.2
50	B3	48	MET	2.2
39	BR	68	ARG	2.2
7	CH	112	ASP	2.2
30	BH	129	GLU	2.2
31	BJ	123	LYS	2.2
36	BO	99	TYR	2.2
39	DR	81	LYS	2.2
52	DI	115	ASP	2.2
36	BO	116	GLN	2.2
39	DR	6	GLN	2.2
52	BI	114	ALA	2.2
52	DI	103	ALA	2.2
2	AC	156	LEU	2.2
7	AH	31	LEU	2.2
20	AB	56	LEU	2.2
30	DH	54	LEU	2.2
3	AD	106	PHE	2.2
3	AD	60	VAL	2.2
8	CI	110	VAL	2.2
14	AO	51	SER	2.2
18	AS	61	VAL	2.2
21	CU	6	ARG	2.2
24	BV	30	ILE	2.2
26	BD	46	ARG	2.2
30	DH	131	SER	2.2
35	DN	69	ARG	2.2
4	AE	9	GLU	2.2
7	AH	101	ALA	2.2
24	BV	94	ALA	2.2
35	BN	73	ASN	2.2
19	CT	23	ARG	2.2
36	BO	29	HIS	2.2
6	AG	90	VAL	2.2
27	DE	119	ILE	2.2
15	CP	68	SER	2.2
52	BI	98	GLY	2.2
33	BL	72	ALA	2.2
2	CC	156	LEU	2.2
6	AG	15	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
21	AU	40	PRO	2.2
52	BI	22	PRO	2.2
26	BD	112	THR	2.2
29	BG	33	THR	2.2
35	BN	120	GLU	2.2
38	DQ	88	GLU	2.2
26	DD	72	GLY	2.2
2	AC	167	TYR	2.2
8	CI	63	TYR	2.2
31	BJ	133	ALA	2.2
38	DQ	31	TYR	2.2
4	AE	10	LEU	2.2
25	DC	166	ARG	2.2
26	DD	103	ASP	2.2
33	BL	104	GLN	2.2
51	B4	36	ARG	2.2
19	CT	30	PHE	2.2
41	BT	51	PHE	2.2
25	BC	266	ILE	2.2
41	DT	56	GLU	2.2
46	DZ	11	GLU	2.2
12	CM	2	ARG	2.2
13	CN	58	ARG	2.2
25	BC	204	LEU	2.2
28	BF	117	SER	2.2
33	BL	33	ARG	2.2
36	BO	18	LEU	2.2
52	DI	64	ARG	2.2
43	BW	63	ASP	2.2
26	BD	165	MET	2.2
34	DM	47	GLU	2.2
37	BP	2	ASN	2.2
1	AA	880	C	2.2
11	CL	68	GLY	2.2
8	AI	123	ARG	2.2
3	AD	121	ALA	2.2
35	BN	79	LEU	2.2
46	BZ	32	LEU	2.2
26	BD	145	SER	2.2
36	DO	5	SER	2.2
37	BP	35	SER	2.2
47	D0	10	SER	2.2

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Mol	Chain	Res	Type	RSRZ
12	AM	99	GLN	2.2
27	DE	175	ILE	2.2
40	DS	4	ILE	2.2
42	BU	82	VAL	2.2
43	BW	36	ILE	2.2
12	CM	110	GLY	2.2
30	DH	96	THR	2.2
47	B0	46	GLY	2.2
2	CC	200	TRP	2.1
29	BG	106	LEU	2.1
6	AG	85	GLN	2.1
20	AB	109	SER	2.1
23	BB	846	U	2.1
27	DE	90	GLN	2.1
33	DL	54	GLN	2.1
43	DW	80	SER	2.1
2	AC	180	ASP	2.1
4	CE	63	MET	2.1
9	AJ	30	LYS	2.1
10	CK	17	ASP	2.1
20	AB	39	ILE	2.1
26	BD	13	ARG	2.1
32	BK	84	CYS	2.1
38	BQ	49	ARG	2.1
40	BS	92	ARG	2.1
43	DW	40	ARG	2.1
45	BY	29	ARG	2.1
45	DY	53	MET	2.1
34	DM	37	GLY	2.1
6	AG	102	TRP	2.1
7	AH	37	ASN	2.1
20	CB	160	LEU	2.1
27	BE	165	HIS	2.1
45	DY	33	HIS	2.1
19	CT	10	ALA	2.1
30	BH	59	ALA	2.1
6	CG	24	LYS	2.1
7	AH	45	ILE	2.1
11	AL	24	GLU	2.1
36	DO	13	ARG	2.1
36	DO	20	GLU	2.1
41	BT	33	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
41	DT	43	ILE	2.1
42	DU	41	VAL	2.1
33	BL	112	LEU	2.1
32	BK	104	THR	2.1
35	BN	21	PHE	2.1
43	BW	62	ALA	2.1
23	BB	2152	G	2.1
52	BI	37	PHE	2.1
25	BC	211	ARG	2.1
30	DH	41	LYS	2.1
32	DK	71	ARG	2.1
2	AC	152	VAL	2.1
16	AQ	58	VAL	2.1
31	BJ	73	VAL	2.1
33	BL	116	VAL	2.1
38	DQ	19	GLN	2.1
26	BD	93	GLY	2.1
26	BD	198	GLY	2.1
41	BT	90	GLY	2.1
42	DU	47	PRO	2.1
9	AJ	73	LEU	2.1
15	CP	54	LEU	2.1
11	CL	19	ASN	2.1
20	CB	159	ALA	2.1
29	BG	50	THR	2.1
35	BN	25	ALA	2.1
48	D1	51	ALA	2.1
13	AN	29	ILE	2.1
29	BG	41	GLU	2.1
41	BT	30	ILE	2.1
50	B3	3	ILE	2.1
52	DI	34	ILE	2.1
1	CA	412	A	2.1
2	AC	133	MET	2.1
5	CF	62	MET	2.1
8	CI	9	GLY	2.1
9	AJ	71	LEU	2.1
11	CL	90	PRO	2.1
21	AU	10	PRO	2.1
35	DN	98	LEU	2.1
50	D3	56	LEU	2.1
8	CI	118	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
23	BB	1174	U	2.1
29	BG	38	ASP	2.1
27	BE	86	ALA	2.1
36	BO	70	ALA	2.1
41	BT	83	ALA	2.1
41	DT	77	ARG	2.1
5	CF	92	THR	2.1
6	AG	73	GLU	2.1
13	AN	34	ASN	2.1
25	DC	171	VAL	2.1
35	BN	47	VAL	2.1
42	DU	92	VAL	2.1
14	AO	65	LEU	2.1
24	DV	61	LEU	2.1
25	DC	224	MET	2.1
49	D2	22	MET	2.1
26	DD	152	PRO	2.1
50	B3	45	PRO	2.1
51	B4	31	PRO	2.1
21	CU	44	ARG	2.1
25	DC	261	ARG	2.1
45	BY	10	ARG	2.1
6	CG	112	ASP	2.1
20	CB	33	ALA	2.1
27	DE	201	ALA	2.1
29	DG	57	TYR	2.1
29	BG	114	HIS	2.1
32	BK	89	ASN	2.1
39	BR	51	VAL	2.1
43	BW	56	HIS	2.1
25	DC	238	ASN	2.1
48	D1	28	THR	2.1
17	AR	64	LEU	2.1
34	DM	102	LEU	2.1
25	DC	187	CYS	2.1
33	BL	58	TYR	2.1
1	AA	845	A	2.1
10	AK	73	VAL	2.1
23	DB	2602	A	2.1
24	BV	72	VAL	2.1
25	DC	136	VAL	2.1
30	BH	114	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
23	DB	2833	U	2.1
24	DV	4	ILE	2.1
28	DF	84	ILE	2.1
33	DL	142	ILE	2.1
43	DW	36	ILE	2.1
3	AD	104	MET	2.1
20	AB	41	ASN	2.1
27	DE	95	LYS	2.1
28	BF	75	GLY	2.1
38	BQ	36	GLN	2.1
39	BR	90	ARG	2.1
46	DZ	49	ARG	2.1
15	CP	15	PRO	2.1
2	AC	160	GLU	2.1
11	CL	12	ALA	2.1
50	B3	10	ALA	2.1
52	BI	109	ALA	2.1
3	AD	82	LYS	2.1
8	CI	54	VAL	2.1
13	AN	33	VAL	2.1
9	AJ	56	HIS	2.1
19	CT	3	ILE	2.1
32	BK	35	VAL	2.1
34	BM	131	VAL	2.1
50	B3	6	VAL	2.1
52	BI	120	ASP	2.1
29	DG	36	LEU	2.1
31	BJ	82	GLY	2.1
44	BX	36	GLN	2.1
25	BC	251	THR	2.1
32	BK	79	PHE	2.1
20	AB	200	PRO	2.1
6	CG	42	VAL	2.1
11	AL	118	VAL	2.1
16	CQ	29	LYS	2.1
27	DE	87	ALA	2.1
28	BF	57	ALA	2.1
31	BJ	85	LYS	2.1
44	BX	33	ALA	2.1
26	BD	5	VAL	2.1
4	CE	118	GLY	2.1
26	DD	43	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
29	DG	46	ASP	2.1
40	DS	46	LEU	2.1
41	BT	11	LEU	2.1
44	BX	6	LEU	2.1
46	DZ	35	ASP	2.1
8	CI	49	GLN	2.1
26	BD	49	GLN	2.1
46	DZ	60	PHE	2.1
6	CG	79	VAL	2.1
37	DP	87	ARG	2.1
42	DU	35	VAL	2.1
7	AH	73	SER	2.1
8	AI	71	ILE	2.1
33	DL	105	ILE	2.1
33	DL	135	ILE	2.1
50	B3	58	ILE	2.1
29	BG	13	GLY	2.1
52	DI	118	GLY	2.1
19	AT	50	PHE	2.1
5	CF	59	TYR	2.1
9	AJ	37	ARG	2.1
23	DB	100	U	2.1
26	DD	81	GLU	2.1
8	AI	66	VAL	2.1
10	CK	116	PRO	2.1
9	AJ	25	ILE	2.1
28	BF	103	ILE	2.1
6	AG	14	ASP	2.1
17	CR	21	ASP	2.1
37	DP	88	ARG	2.0
41	BT	66	LYS	2.0
1	AA	842	U	2.0
9	AJ	44	THR	2.0
23	BB	1176	U	2.0
27	DE	163	ASN	2.0
30	DH	144	VAL	2.0
35	BN	10	LEU	2.0
41	DT	14	PRO	2.0
42	DU	10	VAL	2.0
52	BI	56	VAL	2.0
6	CG	7	GLY	2.0
7	AH	60	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
43	BW	41	GLY	2.0
4	AE	110	MET	2.0
20	CB	125	PHE	2.0
27	DE	41	GLN	2.0
29	BG	127	GLN	2.0
49	D2	19	ARG	2.0
50	B3	29	ARG	2.0
3	AD	52	VAL	2.0
8	AI	18	VAL	2.0
26	DD	9	VAL	2.0
52	DI	113	ALA	2.0
37	DP	109	ILE	2.0
38	DQ	59	LEU	2.0
11	AL	43	LYS	2.0
52	DI	28	GLY	2.0
3	AD	177	MET	2.0
25	BC	29	PHE	2.0
28	BF	149	ARG	2.0
16	AQ	62	GLU	2.0
18	CS	63	ASP	2.0
27	DE	198	GLU	2.0
36	DO	34	HIS	2.0
13	CN	44	VAL	2.0
34	BM	63	ILE	2.0
38	BQ	117	ALA	2.0
42	BU	27	VAL	2.0
6	CG	11	ILE	2.0
12	AM	3	ILE	2.0
17	CR	66	LEU	2.0
25	DC	103	ILE	2.0
33	BL	22	GLY	2.0
47	D0	52	LYS	2.0
51	B4	18	LYS	2.0
33	DL	30	THR	2.0
33	DL	69	ARG	2.0
44	DX	29	ARG	2.0
23	BB	2180	U	2.0
25	DC	114	GLN	2.0
37	DP	111	GLU	2.0
25	BC	52	HIS	2.0
5	CF	74	LEU	2.0
8	CI	114	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
19	CT	15	LYS	2.0
34	BM	93	VAL	2.0
1	AA	74	A	2.0
10	AK	92	ARG	2.0
19	AT	69	ASN	2.0
26	BD	129	THR	2.0
30	BH	11	ASN	2.0
9	AJ	47	GLU	2.0
5	AF	41	ASP	2.0
5	CF	4	TYR	2.0
6	AG	86	VAL	2.0
13	AN	32	ASP	2.0
20	AB	17	HIS	2.0
26	DD	134	HIS	2.0
31	BJ	125	TYR	2.0
52	BI	71	LYS	2.0
19	AT	38	ILE	2.0
20	AB	94	ARG	2.0
26	DD	54	ALA	2.0
37	DP	96	LEU	2.0
26	BD	87	GLY	2.0
21	AU	36	PHE	2.0
23	DB	1068	G	2.0
44	DX	62	GLY	2.0
20	AB	24	PRO	2.0
20	AB	28	PRO	2.0
3	AD	178	GLU	2.0
5	CF	88	MET	2.0
1	CA	999	C	2.0
13	AN	42	ASN	2.0
23	BB	1057	A	2.0
26	DD	136	ASN	2.0
29	DG	128	THR	2.0
33	BL	106	GLU	2.0
31	BJ	86	GLN	2.0
3	AD	129	VAL	2.0
29	BG	22	VAL	2.0
30	BH	68	ARG	2.0
30	BH	103	VAL	2.0
32	BK	8	LEU	2.0
33	DL	57	LEU	2.0
34	BM	89	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
37	BP	91	VAL	2.0
51	B4	17	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
54	MG	DB	3082	1/1	0.82	0.29	6.31	61,61,61,61	0
54	MG	BB	3086	1/1	0.95	0.29	5.28	94,94,94,94	0
53	KSG	CA	1601	26/26	0.83	0.33	3.77	53,53,53,53	0
53	KSG	AA	1601	26/26	0.83	0.30	2.61	53,53,53,53	0
54	MG	BB	3100	1/1	0.73	0.22	2.12	37,37,37,37	1
54	MG	DB	3025	1/1	0.91	0.26	1.72	20,20,20,20	0
54	MG	BB	3078	1/1	0.95	0.28	1.33	49,49,49,49	0
54	MG	AA	1631	1/1	0.68	0.26	1.07	143,143,143,143	0
54	MG	DB	3086	1/1	0.94	0.24	0.94	25,25,25,25	0
54	MG	DB	3094	1/1	0.89	0.21	0.62	33,33,33,33	0
54	MG	BB	3110	1/1	0.92	0.22	0.58	79,79,79,79	0
54	MG	DN	201	1/1	0.94	0.26	0.13	38,38,38,38	0
54	MG	DB	3067	1/1	0.91	0.28	-0.22	11,11,11,11	0
54	MG	BB	3081	1/1	0.81	0.21	-0.26	34,34,34,34	0
54	MG	BB	3012	1/1	0.97	0.19	-0.35	75,75,75,75	0
54	MG	AA	1608	1/1	0.66	0.19	-0.49	57,57,57,57	0
54	MG	DB	3085	1/1	0.86	0.21	-0.49	64,64,64,64	0
54	MG	AA	1616	1/1	0.71	0.20	-0.58	86,86,86,86	0
54	MG	BB	3034	1/1	0.67	0.23	-0.62	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	BB	3018	1/1	0.75	0.22	-0.72	42,42,42,42	0
54	MG	BB	3073	1/1	0.92	0.21	-0.80	55,55,55,55	0
54	MG	BB	3048	1/1	0.84	0.14	-0.83	39,39,39,39	0
54	MG	DB	3050	1/1	0.92	0.20	-0.86	42,42,42,42	0
54	MG	DB	3109	1/1	0.92	0.19	-0.86	57,57,57,57	0
54	MG	DB	3056	1/1	0.90	0.12	-0.88	67,67,67,67	0
54	MG	DB	3062	1/1	0.80	0.12	-1.08	24,24,24,24	0
54	MG	BB	3005	1/1	0.85	0.15	-1.08	13,13,13,13	0
54	MG	BB	3062	1/1	0.92	0.21	-1.12	34,34,34,34	0
54	MG	AA	1602	1/1	0.83	0.19	-1.18	34,34,34,34	0
54	MG	DB	3055	1/1	0.93	0.20	-1.19	20,20,20,20	0
54	MG	DB	3012	1/1	0.91	0.17	-1.20	6,6,6,6	0
54	MG	AA	1653	1/1	0.92	0.15	-1.21	30,30,30,30	0
54	MG	DB	3003	1/1	0.87	0.13	-1.27	8,8,8,8	0
54	MG	CA	1615	1/1	0.95	0.14	-1.36	54,54,54,54	0
54	MG	BB	3092	1/1	0.94	0.16	-1.43	56,56,56,56	0
54	MG	AA	1610	1/1	0.98	0.16	-1.44	10,10,10,10	0
54	MG	BB	3088	1/1	0.90	0.16	-1.45	25,25,25,25	0
54	MG	DB	3024	1/1	0.85	0.12	-1.52	27,27,27,27	0
54	MG	DB	3029	1/1	0.85	0.19	-1.56	22,22,22,22	0
54	MG	BB	3098	1/1	0.94	0.15	-1.63	49,49,49,49	0
54	MG	DB	3076	1/1	0.89	0.15	-1.68	69,69,69,69	0
54	MG	DB	3027	1/1	0.88	0.15	-1.75	37,37,37,37	0
54	MG	DB	3046	1/1	0.95	0.18	-1.76	18,18,18,18	0
54	MG	CA	1640	1/1	0.80	0.12	-1.86	139,139,139,139	0
54	MG	DB	3001	1/1	0.93	0.17	-1.89	5,5,5,5	0
54	MG	CA	1642	1/1	0.91	0.13	-1.93	70,70,70,70	0
54	MG	BB	3029	1/1	0.97	0.10	-1.97	7,7,7,7	0
54	MG	CA	1613	1/1	0.94	0.10	-2.14	63,63,63,63	0
54	MG	DB	3108	1/1	0.75	0.12	-2.15	24,24,24,24	0
54	MG	DB	3092	1/1	0.95	0.06	-2.25	16,16,16,16	0
54	MG	CA	1617	1/1	0.92	0.07	-2.29	45,45,45,45	0
54	MG	BB	3065	1/1	0.84	0.14	-2.33	47,47,47,47	0
54	MG	BB	3096	1/1	0.90	0.12	-2.37	39,39,39,39	0
54	MG	BB	3040	1/1	0.95	0.14	-2.44	18,18,18,18	0
54	MG	BB	3103	1/1	0.96	0.14	-2.50	11,11,11,11	0
54	MG	BB	3087	1/1	0.92	0.10	-2.53	53,53,53,53	0
54	MG	BB	3021	1/1	0.82	0.17	-2.70	24,24,24,24	0
54	MG	BB	3094	1/1	0.83	0.14	-2.84	16,16,16,16	0
54	MG	BB	3051	1/1	0.86	0.10	-2.87	74,74,74,74	0
54	MG	DB	3010	1/1	0.93	0.08	-2.88	8,8,8,8	0
54	MG	DB	3014	1/1	0.94	0.14	-2.88	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
54	MG	AA	1661	1/1	0.79	0.05	-2.90	83,83,83,83	0
54	MG	DB	3022	1/1	0.93	0.07	-2.92	5,5,5,5	0
54	MG	CA	1602	1/1	0.90	0.12	-2.93	8,8,8,8	0
54	MG	AA	1604	1/1	0.89	0.11	-2.98	25,25,25,25	0
54	MG	DB	3034	1/1	0.92	0.11	-2.99	84,84,84,84	0
54	MG	BB	3056	1/1	0.93	0.14	-3.04	37,37,37,37	0
54	MG	BB	3023	1/1	0.93	0.10	-3.07	27,27,27,27	0
54	MG	BB	3090	1/1	0.89	0.11	-3.12	65,65,65,65	0
54	MG	DB	3072	1/1	0.97	0.10	-3.12	5,5,5,5	0
54	MG	DB	3097	1/1	0.96	0.13	-3.16	19,19,19,19	0
54	MG	CA	1647	1/1	0.91	0.12	-3.26	40,40,40,40	0
54	MG	AA	1654	1/1	0.80	0.11	-3.34	81,81,81,81	0
54	MG	DB	3090	1/1	0.95	0.15	-3.38	106,106,106,106	0
54	MG	CA	1634	1/1	0.94	0.09	-3.52	46,46,46,46	0
54	MG	CA	1606	1/1	0.94	0.08	-3.55	30,30,30,30	0
54	MG	BB	3037	1/1	0.95	0.10	-3.58	21,21,21,21	0
54	MG	AA	1638	1/1	0.95	0.11	-3.59	88,88,88,88	0
54	MG	DB	3077	1/1	0.97	0.13	-3.62	41,41,41,41	0
54	MG	BB	3085	1/1	0.93	0.11	-3.78	67,67,67,67	0
54	MG	AA	1643	1/1	0.97	0.07	-3.80	116,116,116,116	0
54	MG	DB	3068	1/1	0.95	0.12	-3.80	49,49,49,49	0
54	MG	CA	1618	1/1	0.98	0.08	-3.85	26,26,26,26	0
54	MG	CA	1658	1/1	0.90	0.07	-3.91	18,18,18,18	0
54	MG	DB	3106	1/1	0.89	0.12	-4.07	7,7,7,7	0
54	MG	DB	3083	1/1	0.92	0.13	-4.28	18,18,18,18	0
54	MG	BB	3026	1/1	0.91	0.12	-4.30	25,25,25,25	0
54	MG	BB	3011	1/1	0.97	0.12	-4.43	9,9,9,9	0
54	MG	AA	1651	1/1	0.92	0.05	-4.46	64,64,64,64	0
54	MG	BB	3052	1/1	0.90	0.09	-4.49	27,27,27,27	0
54	MG	BB	3083	1/1	0.88	0.14	-4.56	25,25,25,25	0
54	MG	DB	3009	1/1	0.91	0.11	-4.58	10,10,10,10	0
54	MG	DB	3101	1/1	0.95	0.13	-4.78	30,30,30,30	0
54	MG	CA	1635	1/1	0.93	0.11	-4.83	91,91,91,91	0
54	MG	BB	3035	1/1	0.96	0.12	-4.84	25,25,25,25	0
54	MG	BB	3013	1/1	0.91	0.09	-4.84	36,36,36,36	0
54	MG	CA	1657	1/1	0.92	0.13	-4.93	51,51,51,51	0
54	MG	BB	3019	1/1	0.93	0.14	-4.94	22,22,22,22	0
54	MG	AA	1611	1/1	0.95	0.06	-5.03	57,57,57,57	0
54	MG	DB	3007	1/1	0.82	0.14	-5.08	15,15,15,15	0
54	MG	DB	3096	1/1	0.99	0.17	-5.09	31,31,31,31	0
54	MG	DB	3087	1/1	0.96	0.12	-5.12	40,40,40,40	0
54	MG	DB	3026	1/1	0.88	0.14	-5.12	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	BB	3079	1/1	0.94	0.06	-5.43	32,32,32,32	0
54	MG	DB	3075	1/1	0.81	0.14	-5.56	42,42,42,42	0
54	MG	BB	3108	1/1	0.90	0.13	-5.94	22,22,22,22	0
54	MG	BB	3069	1/1	0.96	0.08	-6.00	8,8,8,8	0
54	MG	CA	1648	1/1	0.96	0.06	-6.19	39,39,39,39	0
54	MG	CA	1619	1/1	0.96	0.07	-6.23	9,9,9,9	0
54	MG	DB	3002	1/1	0.91	0.13	-6.44	20,20,20,20	0
54	MG	BB	3061	1/1	0.94	0.06	-6.51	33,33,33,33	0
54	MG	BB	3002	1/1	0.98	0.09	-6.59	9,9,9,9	0
54	MG	DB	3054	1/1	0.97	0.10	-6.68	32,32,32,32	0
54	MG	CA	1645	1/1	0.96	0.08	-7.16	28,28,28,28	0
54	MG	DB	3078	1/1	0.97	0.08	-10.26	5,5,5,5	0
54	MG	BB	3001	1/1	0.85	0.11	-10.88	28,28,28,28	0
54	MG	BB	3004	1/1	0.92	0.08	-	39,39,39,39	0
54	MG	BB	3017	1/1	0.89	0.10	-	54,54,54,54	0
54	MG	DB	3060	1/1	0.83	0.10	-	105,105,105,105	0
54	MG	DB	3035	1/1	0.83	0.16	-	24,24,24,24	0
54	MG	AA	1620	1/1	0.74	0.29	-	116,116,116,116	0
54	MG	DB	3103	1/1	0.82	0.17	-	61,61,61,61	0
54	MG	BB	3076	1/1	0.85	0.12	-	43,43,43,43	0
54	MG	AA	1659	1/1	0.97	0.20	-	157,157,157,157	0
54	MG	CA	1659	1/1	0.91	0.25	-	88,88,88,88	0
54	MG	AA	1625	1/1	0.86	0.40	-	81,81,81,81	0
54	MG	AA	1614	1/1	0.73	0.15	-	67,67,67,67	0
54	MG	CA	1620	1/1	0.96	0.14	-	32,32,32,32	0
54	MG	AA	1627	1/1	0.75	0.33	-	17,17,17,17	1
54	MG	BB	3032	1/1	0.92	0.14	-	31,31,31,31	0
54	MG	DB	3043	1/1	0.94	0.07	-	7,7,7,7	0
54	MG	CA	1637	1/1	0.84	0.15	-	107,107,107,107	0
54	MG	BB	3022	1/1	0.96	0.14	-	55,55,55,55	0
54	MG	AA	1605	1/1	0.92	0.12	-	30,30,30,30	0
54	MG	DB	3005	1/1	0.98	0.09	-	6,6,6,6	0
54	MG	AA	1628	1/1	0.71	0.16	-	63,63,63,63	0
54	MG	DB	3053	1/1	0.88	0.10	-	37,37,37,37	0
54	MG	AA	1609	1/1	0.49	0.20	-	119,119,119,119	0
54	MG	AA	1607	1/1	0.81	0.08	-	82,82,82,82	0
54	MG	BB	3093	1/1	0.92	0.17	-	5,5,5,5	1
54	MG	CA	1646	1/1	0.95	0.05	-	94,94,94,94	0
54	MG	DB	3061	1/1	0.89	0.11	-	48,48,48,48	0
54	MG	DB	3015	1/1	0.87	0.15	-	69,69,69,69	0
54	MG	DB	3006	1/1	0.98	0.08	-	16,16,16,16	0
54	MG	CA	1610	1/1	0.97	0.16	-	61,61,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	AA	1615	1/1	0.30	0.19	-	72,72,72,72	0
54	MG	BB	3084	1/1	0.80	0.11	-	19,19,19,19	0
54	MG	DB	3051	1/1	0.94	0.15	-	57,57,57,57	0
54	MG	AA	1650	1/1	0.44	0.23	-	133,133,133,133	0
54	MG	BB	3014	1/1	0.84	0.17	-	33,33,33,33	0
54	MG	DB	3069	1/1	0.93	0.09	-	50,50,50,50	0
54	MG	AA	1649	1/1	0.85	0.10	-	102,102,102,102	0
54	MG	AA	1660	1/1	0.80	0.10	-	95,95,95,95	0
54	MG	CA	1607	1/1	0.88	0.13	-	126,126,126,126	0
54	MG	CA	1643	1/1	0.95	0.10	-	62,62,62,62	0
54	MG	BB	3024	1/1	0.90	0.17	-	22,22,22,22	0
54	MG	CA	1625	1/1	0.96	0.14	-	24,24,24,24	0
54	MG	CA	1644	1/1	0.82	0.12	-	59,59,59,59	0
54	MG	DB	3020	1/1	0.94	0.24	-	41,41,41,41	0
54	MG	AA	1613	1/1	0.80	0.20	-	108,108,108,108	0
54	MG	CA	1652	1/1	0.94	0.13	-	20,20,20,20	0
54	MG	BB	3007	1/1	0.88	0.14	-	57,57,57,57	0
54	MG	AA	1655	1/1	0.80	0.10	-	67,67,67,67	0
54	MG	DB	3098	1/1	0.88	0.11	-	9,9,9,9	0
54	MG	DB	3032	1/1	0.87	0.11	-	29,29,29,29	0
54	MG	DB	3059	1/1	0.73	0.15	-	75,75,75,75	0
54	MG	BB	3063	1/1	0.87	0.17	-	27,27,27,27	0
54	MG	DB	3042	1/1	0.90	0.16	-	5,5,5,5	0
54	MG	DB	3028	1/1	0.83	0.18	-	29,29,29,29	0
54	MG	BB	3071	1/1	0.91	0.19	-	38,38,38,38	0
54	MG	AA	1640	1/1	0.82	0.10	-	42,42,42,42	0
54	MG	BB	3030	1/1	0.93	0.08	-	45,45,45,45	0
54	MG	CA	1660	1/1	0.96	0.21	-	27,27,27,27	0
54	MG	DB	3021	1/1	0.93	0.12	-	25,25,25,25	0
54	MG	CA	1611	1/1	0.96	0.04	-	43,43,43,43	0
54	MG	BB	3045	1/1	0.97	0.05	-	40,40,40,40	0
54	MG	DB	3066	1/1	0.93	0.15	-	12,12,12,12	0
54	MG	AA	1618	1/1	0.92	0.12	-	102,102,102,102	0
54	MG	BB	3031	1/1	0.86	0.18	-	46,46,46,46	0
54	MG	BB	3080	1/1	0.91	0.09	-	42,42,42,42	0
54	MG	CA	1636	1/1	0.92	0.14	-	13,13,13,13	0
54	MG	AA	1630	1/1	0.95	0.10	-	20,20,20,20	0
54	MG	BB	3060	1/1	0.82	0.16	-	81,81,81,81	0
54	MG	CA	1631	1/1	0.58	0.25	-	21,21,21,21	1
54	MG	BB	3049	1/1	0.89	0.15	-	12,12,12,12	0
54	MG	CA	1654	1/1	0.97	0.19	-	32,32,32,32	0
54	MG	DB	3037	1/1	0.94	0.10	-	8,8,8,8	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	BB	3059	1/1	0.90	0.10	-	16,16,16,16	0
54	MG	AA	1621	1/1	0.73	0.14	-	96,96,96,96	0
54	MG	DB	3099	1/1	0.96	0.17	-	11,11,11,11	0
54	MG	BB	3082	1/1	0.64	0.39	-	63,63,63,63	0
54	MG	CA	1629	1/1	0.62	0.60	-	62,62,62,62	1
54	MG	AA	1633	1/1	0.82	0.15	-	92,92,92,92	0
54	MG	BB	3020	1/1	0.96	0.07	-	38,38,38,38	0
54	MG	CA	1626	1/1	0.95	0.11	-	17,17,17,17	0
54	MG	BB	3016	1/1	0.91	0.18	-	19,19,19,19	0
54	MG	CA	1653	1/1	0.95	0.07	-	48,48,48,48	0
54	MG	BB	3010	1/1	0.89	0.16	-	74,74,74,74	0
54	MG	DB	3052	1/1	0.78	0.25	-	42,42,42,42	0
54	MG	DB	3079	1/1	0.93	0.14	-	7,7,7,7	0
54	MG	AA	1657	1/1	0.81	0.27	-	83,83,83,83	0
54	MG	AA	1622	1/1	0.96	0.06	-	16,16,16,16	0
54	MG	AA	1612	1/1	0.80	0.12	-	52,52,52,52	0
54	MG	DB	3019	1/1	0.99	0.03	-	8,8,8,8	0
54	MG	CA	1656	1/1	0.97	0.10	-	60,60,60,60	0
54	MG	CA	1603	1/1	0.94	0.12	-	32,32,32,32	0
54	MG	DB	3045	1/1	0.95	0.07	-	45,45,45,45	0
54	MG	DB	3033	1/1	0.77	0.26	-	88,88,88,88	0
54	MG	CA	1633	1/1	0.96	0.10	-	50,50,50,50	0
54	MG	DB	3071	1/1	0.72	0.13	-	27,27,27,27	0
54	MG	DB	3089	1/1	0.97	0.28	-	28,28,28,28	0
54	MG	BB	3106	1/1	0.92	0.25	-	62,62,62,62	0
54	MG	CA	1605	1/1	0.96	0.09	-	37,37,37,37	0
54	MG	DB	3039	1/1	0.93	0.11	-	5,5,5,5	0
54	MG	BB	3075	1/1	0.91	0.25	-	35,35,35,35	0
54	MG	DB	3084	1/1	0.97	0.17	-	18,18,18,18	0
54	MG	CA	1616	1/1	0.59	0.26	-	153,153,153,153	0
54	MG	BB	3003	1/1	0.94	0.11	-	23,23,23,23	0
54	MG	AA	1635	1/1	0.97	0.07	-	75,75,75,75	0
54	MG	CA	1638	1/1	0.88	0.21	-	60,60,60,60	0
54	MG	BB	3097	1/1	0.72	0.13	-	96,96,96,96	0
54	MG	BB	3041	1/1	0.81	0.18	-	47,47,47,47	0
54	MG	DB	3100	1/1	0.94	0.14	-	6,6,6,6	0
54	MG	AA	1606	1/1	0.91	0.06	-	28,28,28,28	0
54	MG	BB	3008	1/1	0.74	0.16	-	82,82,82,82	0
54	MG	BB	3095	1/1	0.98	0.06	-	48,48,48,48	0
54	MG	DB	3104	1/1	0.94	0.19	-	44,44,44,44	0
54	MG	CA	1628	1/1	0.85	0.33	-	38,38,38,38	1
54	MG	CA	1604	1/1	0.95	0.15	-	49,49,49,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	CA	1621	1/1	0.92	0.09	-	46,46,46,46	0
54	MG	BB	3058	1/1	0.97	0.12	-	33,33,33,33	0
54	MG	DB	3017	1/1	0.74	0.23	-	22,22,22,22	0
54	MG	BB	3089	1/1	0.93	0.23	-	82,82,82,82	0
54	MG	AA	1636	1/1	0.94	0.14	-	83,83,83,83	0
54	MG	BB	3074	1/1	0.87	0.12	-	7,7,7,7	0
54	MG	DB	3095	1/1	0.91	0.09	-	44,44,44,44	0
54	MG	AA	1644	1/1	0.95	0.17	-	81,81,81,81	0
54	MG	BB	3099	1/1	0.89	0.13	-	58,58,58,58	0
54	MG	BB	3064	1/1	0.88	0.16	-	71,71,71,71	0
54	MG	AA	1647	1/1	0.87	0.60	-	84,84,84,84	0
54	MG	CA	1632	1/1	0.94	0.12	-	34,34,34,34	0
54	MG	AA	1656	1/1	0.86	0.14	-	100,100,100,100	0
54	MG	AA	1642	1/1	0.83	0.15	-	107,107,107,107	0
54	MG	AA	1617	1/1	0.87	0.13	-	65,65,65,65	0
54	MG	CA	1627	1/1	0.85	0.19	-	114,114,114,114	0
54	MG	BB	3036	1/1	0.89	0.10	-	29,29,29,29	0
54	MG	CA	1624	1/1	0.32	0.29	-	162,162,162,162	0
54	MG	CA	1630	1/1	0.86	0.09	-	49,49,49,49	0
54	MG	BB	3091	1/1	0.89	0.16	-	42,42,42,42	0
54	MG	BB	3015	1/1	0.97	0.10	-	23,23,23,23	0
54	MG	DB	3018	1/1	0.92	0.12	-	39,39,39,39	0
54	MG	AA	1624	1/1	0.88	0.12	-	25,25,25,25	1
54	MG	AA	1658	1/1	0.80	0.12	-	113,113,113,113	0
54	MG	BB	3033	1/1	0.75	0.25	-	113,113,113,113	0
54	MG	DB	3038	1/1	0.94	0.08	-	33,33,33,33	0
54	MG	DB	3031	1/1	0.97	0.10	-	40,40,40,40	0
54	MG	AA	1652	1/1	0.88	0.14	-	105,105,105,105	0
54	MG	BB	3006	1/1	0.95	0.08	-	20,20,20,20	0
54	MG	DB	3011	1/1	0.94	0.17	-	23,23,23,23	0
54	MG	CA	1609	1/1	0.78	0.19	-	132,132,132,132	0
54	MG	BB	3039	1/1	0.91	0.13	-	26,26,26,26	0
54	MG	DB	3080	1/1	0.89	0.12	-	34,34,34,34	0
54	MG	BB	3101	1/1	0.89	0.10	-	14,14,14,14	0
54	MG	DB	3036	1/1	0.94	0.17	-	25,25,25,25	0
54	MG	DB	3081	1/1	0.91	0.15	-	110,110,110,110	0
54	MG	DB	3065	1/1	0.95	0.16	-	25,25,25,25	0
54	MG	DB	3093	1/1	0.86	0.76	-	48,48,48,48	1
54	MG	BB	3046	1/1	0.82	0.17	-	63,63,63,63	0
54	MG	AA	1639	1/1	0.46	0.19	-	141,141,141,141	0
54	MG	AA	1603	1/1	0.80	0.14	-	74,74,74,74	0
54	MG	DB	3057	1/1	0.81	0.20	-	56,56,56,56	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	DB	3040	1/1	0.85	0.20	-	28,28,28,28	0
54	MG	BB	3047	1/1	0.69	0.29	-	128,128,128,128	0
54	MG	BB	3067	1/1	0.85	0.16	-	56,56,56,56	0
54	MG	DB	3088	1/1	0.96	0.21	-	51,51,51,51	0
54	MG	DB	3070	1/1	0.94	0.07	-	60,60,60,60	0
54	MG	BB	3077	1/1	0.74	0.16	-	37,37,37,37	0
54	MG	BB	3043	1/1	0.88	0.24	-	63,63,63,63	0
54	MG	DB	3049	1/1	0.78	0.10	-	79,79,79,79	0
54	MG	BB	3038	1/1	0.91	0.13	-	82,82,82,82	0
54	MG	CA	1649	1/1	0.94	0.31	-	86,86,86,86	0
54	MG	CA	1623	1/1	0.50	0.14	-	77,77,77,77	0
54	MG	BB	3050	1/1	0.95	0.13	-	20,20,20,20	0
54	MG	DB	3073	1/1	0.94	0.11	-	44,44,44,44	0
54	MG	CA	1662	1/1	0.97	0.11	-	38,38,38,38	0
54	MG	BB	3044	1/1	0.95	0.07	-	39,39,39,39	0
54	MG	BB	3107	1/1	0.82	0.13	-	37,37,37,37	0
54	MG	BB	3057	1/1	0.86	0.18	-	19,19,19,19	0
54	MG	CA	1608	1/1	0.87	0.06	-	31,31,31,31	0
54	MG	AA	1629	1/1	0.96	0.10	-	62,62,62,62	0
54	MG	BB	3028	1/1	0.80	0.22	-	66,66,66,66	0
54	MG	DB	3048	1/1	0.95	0.07	-	7,7,7,7	0
54	MG	CA	1655	1/1	0.95	0.05	-	26,26,26,26	0
54	MG	DB	3102	1/1	0.95	0.11	-	13,13,13,13	0
54	MG	BB	3053	1/1	0.87	0.12	-	36,36,36,36	0
54	MG	BB	3070	1/1	0.98	0.09	-	55,55,55,55	0
54	MG	AA	1648	1/1	0.97	0.12	-	31,31,31,31	0
54	MG	AA	1634	1/1	0.90	0.19	-	73,73,73,73	0
54	MG	DB	3023	1/1	0.78	0.08	-	44,44,44,44	0
54	MG	DB	3016	1/1	0.80	0.14	-	25,25,25,25	0
54	MG	BB	3072	1/1	0.88	0.12	-	26,26,26,26	0
54	MG	AA	1632	1/1	0.93	0.07	-	44,44,44,44	0
54	MG	CA	1641	1/1	0.94	0.08	-	40,40,40,40	0
54	MG	AA	1619	1/1	0.74	0.13	-	33,33,33,33	0
54	MG	DB	3041	1/1	0.97	0.13	-	29,29,29,29	0
54	MG	CA	1650	1/1	0.94	0.08	-	67,67,67,67	0
54	MG	BB	3054	1/1	0.92	0.11	-	77,77,77,77	0
54	MG	CA	1622	1/1	0.63	0.34	-	132,132,132,132	0
54	MG	BB	3042	1/1	0.87	0.07	-	74,74,74,74	0
54	MG	DB	3091	1/1	0.97	0.16	-	39,39,39,39	0
54	MG	BB	3009	1/1	0.95	0.08	-	35,35,35,35	0
54	MG	DB	3063	1/1	0.98	0.11	-	42,42,42,42	0
54	MG	DB	3044	1/1	0.80	0.09	-	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
54	MG	DB	3058	1/1	0.85	0.20	-	132,132,132,132	0
54	MG	CA	1651	1/1	0.84	0.16	-	122,122,122,122	0
54	MG	AA	1641	1/1	0.95	0.09	-	72,72,72,72	0
54	MG	DB	3105	1/1	0.94	0.08	-	24,24,24,24	0
54	MG	CA	1639	1/1	0.93	0.04	-	41,41,41,41	0
54	MG	DB	3008	1/1	0.96	0.19	-	39,39,39,39	0
54	MG	AA	1646	1/1	0.92	0.15	-	103,103,103,103	0
54	MG	DB	3047	1/1	0.89	0.09	-	29,29,29,29	0
54	MG	BB	3025	1/1	0.97	0.16	-	67,67,67,67	0
54	MG	DB	3064	1/1	0.77	0.27	-	66,66,66,66	1
54	MG	AA	1623	1/1	0.88	0.32	-	143,143,143,143	0
54	MG	AA	1626	1/1	0.61	0.43	-	82,82,82,82	1
54	MG	BB	3104	1/1	0.88	0.13	-	33,33,33,33	0
54	MG	CA	1663	1/1	0.92	0.15	-	20,20,20,20	0
54	MG	BB	3055	1/1	0.95	0.12	-	19,19,19,19	0
54	MG	AA	1645	1/1	0.89	0.16	-	97,97,97,97	0
54	MG	CA	1661	1/1	0.89	0.29	-	67,67,67,67	0
54	MG	DB	3013	1/1	0.67	0.20	-	32,32,32,32	0
54	MG	BB	3068	1/1	0.92	0.12	-	39,39,39,39	0
54	MG	DB	3074	1/1	0.97	0.13	-	12,12,12,12	0
54	MG	AA	1637	1/1	0.67	0.62	-	115,115,115,115	0
54	MG	CA	1614	1/1	0.93	0.13	-	97,97,97,97	0
54	MG	BB	3027	1/1	0.93	0.07	-	29,29,29,29	0
54	MG	BB	3105	1/1	0.87	0.15	-	33,33,33,33	0
54	MG	BB	3102	1/1	0.91	0.11	-	34,34,34,34	0
54	MG	DB	3004	1/1	0.94	0.15	-	38,38,38,38	0
54	MG	DB	3107	1/1	0.85	0.13	-	19,19,19,19	0
54	MG	DE	301	1/1	0.80	0.16	-	69,69,69,69	0
54	MG	CA	1612	1/1	0.87	0.10	-	71,71,71,71	0
54	MG	BB	3066	1/1	0.88	0.13	-	42,42,42,42	0
54	MG	DB	3030	1/1	0.92	0.17	-	11,11,11,11	0
54	MG	BB	3109	1/1	0.90	0.15	-	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.