Full wwPDB X-ray Structure Validation Report (i)

Mar 14, 2026 – 09:46 PM UTC

PDB ID	:	4TUA / pdb_00004tua
Title	:	Crystal structure of ASL-Thr bound to Codon ACC-A on the Ribosome
Authors	:	Fagan, C.E.; Dunham, C.M.
Deposited on	:	2014-06-24
Resolution	:	3.60 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

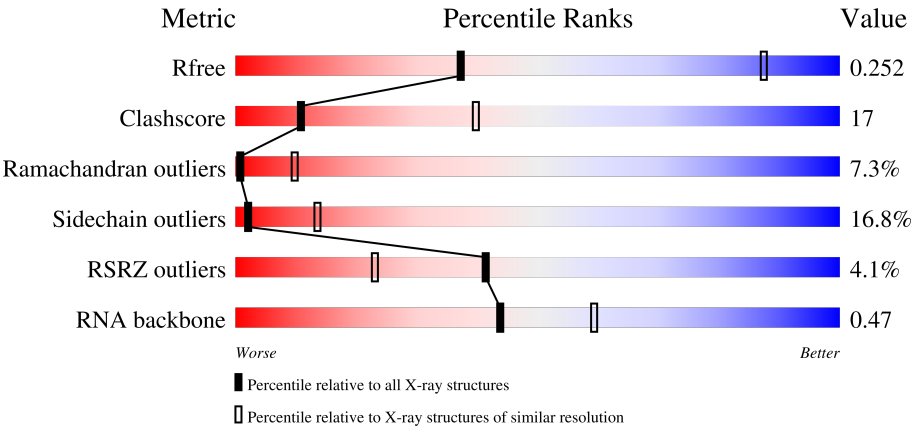
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 3.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)
RNA backbone	3983	1014 (4.14-3.06)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	QA	1522	2% 52% 37% 10% •
1	XA	1522	% 50% 37% 11% •
2	QB	256	7% 46% 38% 8% • 7%
2	XB	256	2% 44% 38% 10% • 7%

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Mol	Chain	Length	Quality of chain
3	QC	239	2% 54% 29% 14%
3	XC	239	3% 55% 26% 14%
4	QD	209	6% 67% 26% 6%
4	XD	209	6% 66% 26% 7%
5	QE	162	3% 55% 30% 7% 7%
5	XE	162	2% 51% 35% 8% 7%
6	QF	101	67% 30%
6	XF	101	62% 33% 5%
7	QG	156	3% 66% 31%
7	XG	156	3% 71% 25%
8	QH	138	4% 57% 38%
8	XH	138	2% 64% 30% 6%
9	QI	128	13% 58% 34% 7%
9	XI	128	13% 54% 32% 12%
10	QJ	105	13% 41% 39% 13% 6%
10	XJ	105	3% 39% 43% 12% 6%
11	QK	129	4% 53% 34% 5% 8%
11	XK	129	3% 61% 29% 8%
12	QL	132	15% 30% 43% 20% 5%
12	XL	132	15% 29% 44% 19% 5%
13	QM	126	10% 48% 37% 11%
13	XM	126	8% 51% 33% 11%
14	QN	61	10% 44% 44% 10%
14	XN	61	10% 49% 38% 11%
15	QO	89	67% 25% 7%

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Mol	Chain	Length	Quality of chain
15	XO	89	66% 25% 8%
16	QP	88	6% 67% 22% 7% 5%
16	XP	88	8% 57% 36% 5%
17	QQ	105	2% 58% 34% 5%
17	XQ	105	5% 68% 23% 5% 5%
18	QR	88	53% 20% 5% 20%
18	XR	88	41% 33% 6% 20%
19	QS	93	4% 38% 33% 16% 10%
19	XS	93	3% 31% 41% 16% 10%
20	QT	106	11% 49% 37% 8% 7%
20	XT	106	15% 38% 43% 12% 7%
21	QU	27	22% 48% 44% 7%
21	XU	27	15% 67% 26% 7%
22	QV	77	57% 34% 9%
22	XV	77	65% 26% 9%
23	QX	25	12% 36% 8% 56%
23	XX	25	16% 32% 12% 56%
24	QY	17	71% 18% 12%
24	XY	17	65% 29% 6%
25	RA	2915	55% 34% 10%
25	YA	2915	56% 34% 9%
26	RB	122	2% 52% 40% 7%
26	YB	122	2% 59% 30% 9%
27	RD	276	4% 52% 36% 10%
27	YD	276	3% 27% 48% 20%

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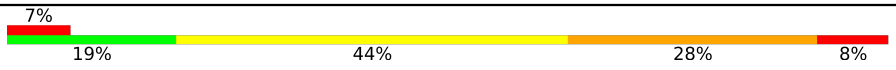
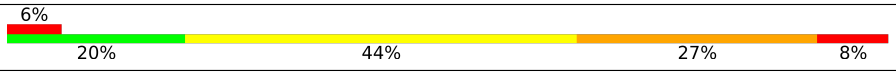
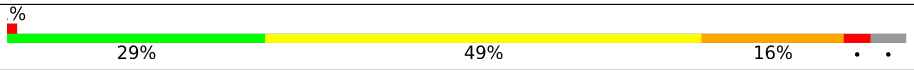
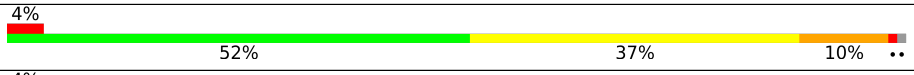
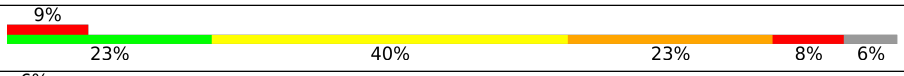
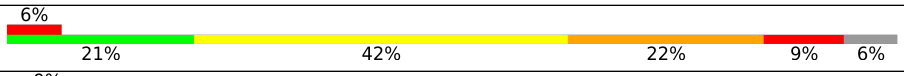
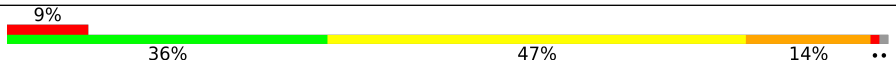
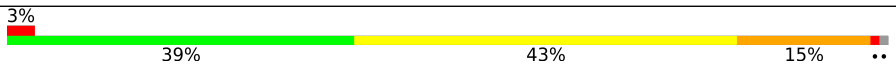
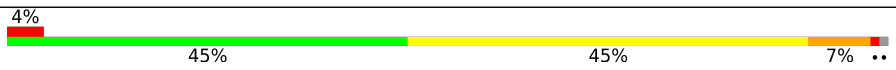
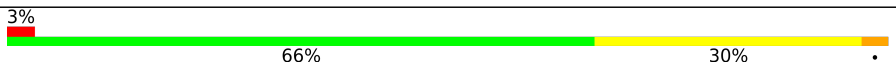
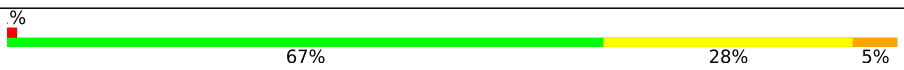
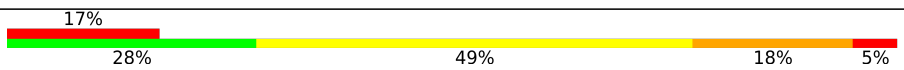
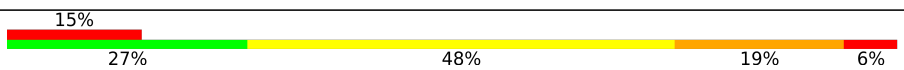
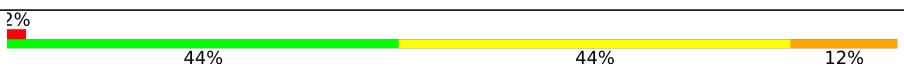
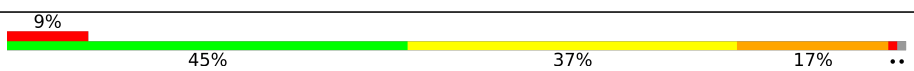
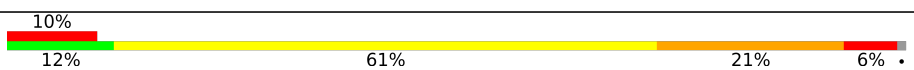
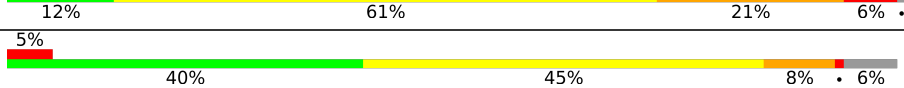
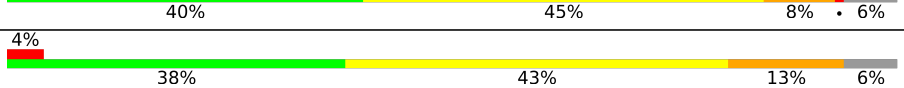
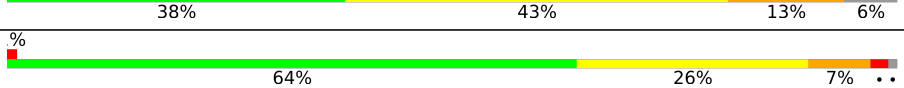
Mol	Chain	Length	Quality of chain
28	RE	206	
28	YE	206	
29	RF	210	
29	YF	210	
30	RG	182	
30	YG	182	
31	RH	180	
31	YH	180	
32	RI	148	
32	YI	148	
33	RN	140	
33	YN	140	
34	RO	122	
34	YO	122	
35	RP	150	
35	YP	150	
36	RQ	141	
36	YQ	141	
37	RR	118	
37	YR	118	
38	RS	112	
38	YS	112	
39	RT	146	
39	YT	146	
40	RU	118	

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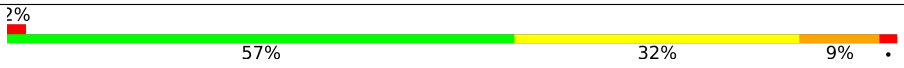
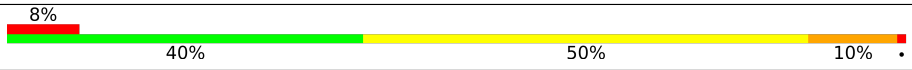
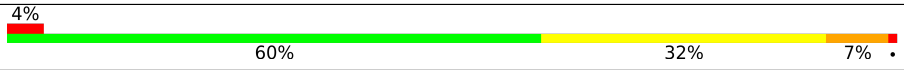
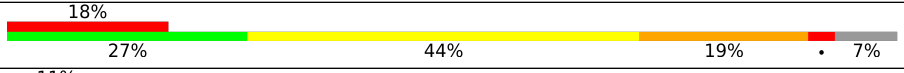
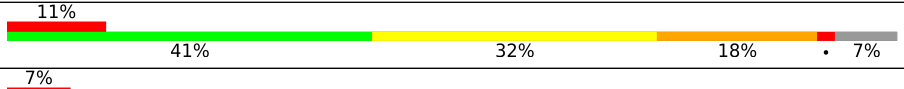
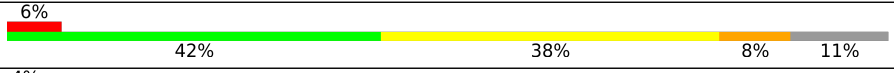
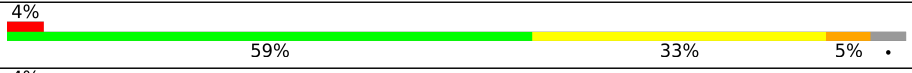
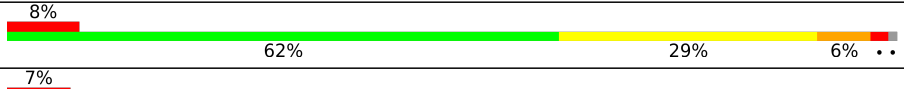
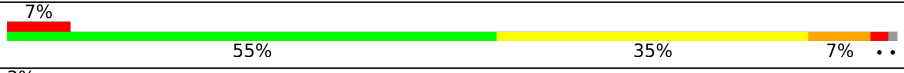
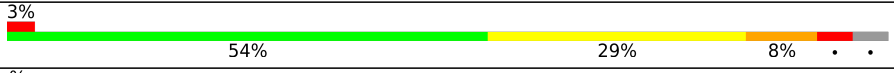
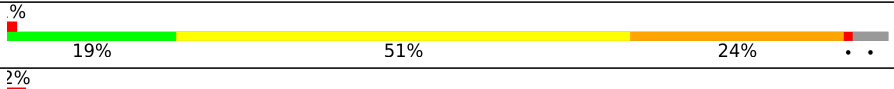
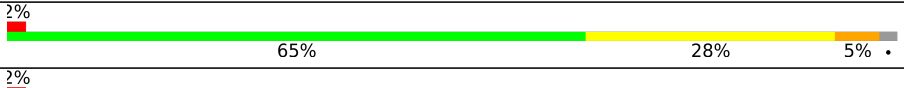
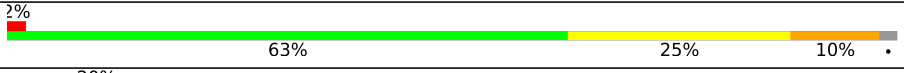
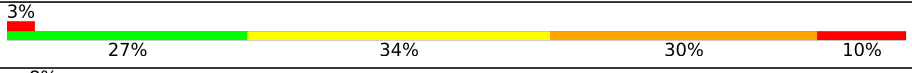
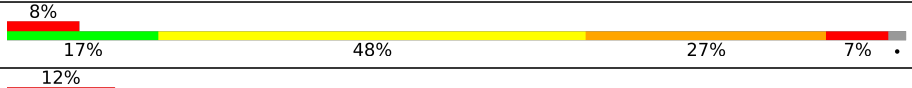
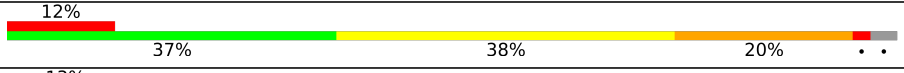
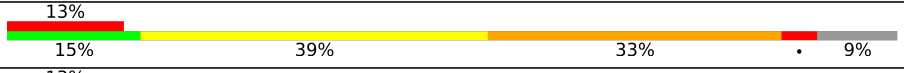
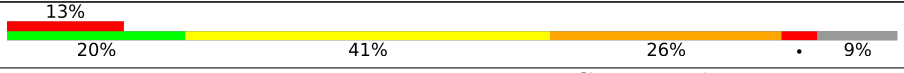
Mol	Chain	Length	Quality of chain
40	YU	118	
41	RV	101	
41	YV	101	
42	RW	113	
42	YW	113	
43	RX	96	
43	YX	96	
44	RY	110	
44	YY	110	
45	RZ	206	
45	YZ	206	
46	R0	85	
46	Y0	85	
47	R1	98	
47	Y1	98	
48	R2	72	
48	Y2	72	
49	R3	60	
49	Y3	60	
50	R4	71	
50	Y4	71	
51	R5	60	
51	Y5	60	
52	R6	54	
52	Y6	54	

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Mol	Chain	Length	Quality of chain
53	R7	49	8% 63% 31% 6%
53	Y7	49	6% 55% 37% 8%
54	R8	65	17% 15% 45% 29% 9%
54	Y8	65	17% 14% 45% 32% 8%
55	R9	37	68% 62% 32% 5%
55	Y9	37	57% 62% 35%
56	Z5	3	67% 33%
56	Z6	3	67% 33%

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
57	MG	RA	3244	-	-	-	X
57	MG	RA	3255	-	-	-	X
57	MG	RA	3265	-	-	-	X
57	MG	YA	3031	-	-	-	X
57	MG	YA	3074	-	-	-	X
57	MG	YA	3280	-	-	-	X

2 Entry composition (i)

There are 59 unique types of molecules in this entry. The entry contains 292311 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	QA	1500	Total	C	N	O	P	0	0	0
			32246	14353	5981	10413	1499			
1	XA	1500	Total	C	N	O	P	0	0	0
			32249	14354	5984	10412	1499			

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	QB	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			
2	XB	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	QC	205	Total	C	N	O	S	0	0	0
			1605	1011	313	280	1			
3	XC	205	Total	C	N	O	S	0	0	0
			1605	1011	313	280	1			

- Molecule 4 is a protein called 30S ribosomal protein S4.

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

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	QE	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			
5	XE	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			

- Molecule 6 is a protein called 30S ribosomal protein S6.

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

- Molecule 7 is a protein called 30S ribosomal protein S7.

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

- Molecule 8 is a protein called 30S ribosomal protein S8.

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

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	QI	127	Total	C	N	O	0	0	0
			1010	639	197	174			
9	XI	127	Total	C	N	O	0	0	0
			1010	639	197	174			

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	QJ	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	XJ	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

- Molecule 11 is a protein called 30S ribosomal protein S11.

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

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	QL	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			
12	XL	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	QM	121	Total	C	N	O	S	0	0	0
			964	597	199	166	2			
13	XM	121	Total	C	N	O	S	0	0	0
			964	597	199	166	2			

- Molecule 14 is a protein called 30S ribosomal protein S14 type Z.

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

- Molecule 15 is a protein called 30S ribosomal protein S15.

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

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	QP	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			
16	XP	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	QQ	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			
17	XQ	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			

- Molecule 18 is a protein called 30S ribosomal protein S18.

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

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	QS	84	Total	C	N	O	S	0	0	0
			674	430	126	116	2			
19	XS	84	Total	C	N	O	S	0	0	0
			674	430	126	116	2			

- Molecule 20 is a protein called 30S ribosomal protein S20.

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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	QU	25	Total	C	N	O	0	0	0
			217	134	52	31			
21	XU	25	Total	C	N	O	0	0	0
			217	134	52	31			

- Molecule 22 is a RNA chain called P-site tRNA-fMet.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	QV	77	Total	C	N	O	P	0	0	0
			1644	732	297	538	77			
22	XV	77	Total	C	N	O	P	0	0	0
			1644	732	297	538	77			

- Molecule 23 is a RNA chain called messenger RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	QX	11	Total	C	N	O	P	0	0	0
			220	98	43	68	11			
23	XX	11	Total	C	N	O	P	0	0	0
			220	98	43	68	11			

- Molecule 24 is a RNA chain called A-site ASL-Thr.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	QY	15	Total	C	N	O	P	0	0	0
			321	143	57	106	15			
24	XY	16	Total	C	N	O	P	0	0	0
			344	153	62	113	16			

- Molecule 25 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	RA	2882	Total	C	N	O	P	0	0	0
			62071	27627	11611	19952	2881			
25	YA	2882	Total	C	N	O	P	0	0	0
			62071	27627	11611	19952	2881			

- Molecule 26 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	RB	120	Total	C	N	O	P	0	0	0
			2573	1146	476	832	119			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	YB	120	Total	C	N	O	P	0	0	0
			2573	1146	476	832	119			

- Molecule 27 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	RD	272	Total	C	N	O	S	0	0	0
			2115	1335	420	357	3			
27	YD	272	Total	C	N	O	S	0	0	0
			2115	1335	420	357	3			

- Molecule 28 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	RE	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			
28	YE	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	RF	202	Total	C	N	O	S	0	0	0
			1585	1011	297	275	2			
29	YF	202	Total	C	N	O	S	0	0	0
			1585	1011	297	275	2			

- Molecule 30 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	RG	181	Total	C	N	O	S	0	0	0
			1474	942	268	260	4			
30	YG	181	Total	C	N	O	S	0	0	0
			1474	942	268	260	4			

- Molecule 31 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	RH	170	Total	C	N	O	S	0	0	0
			1307	829	245	232	1			
31	YH	170	Total	C	N	O	S	0	0	0
			1307	829	245	232	1			

- Molecule 32 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	RI	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			
32	YI	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			

- Molecule 33 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	RN	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			
33	YN	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			

- Molecule 34 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	RO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			
34	YO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			

- Molecule 35 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	RP	150	Total	C	N	O	S	0	0	0
			1145	712	232	198	3			
35	YP	150	Total	C	N	O	S	0	0	0
			1145	712	232	198	3			

- Molecule 36 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	RQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			
36	YQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			

- Molecule 37 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	RR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			
37	YR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			

- Molecule 38 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
38	RS	111	Total	C	N	O	0	0	0
			882	556	176	150			
38	YS	111	Total	C	N	O	0	0	0
			882	556	176	150			

- Molecule 39 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	RT	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			
39	YT	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			

- Molecule 40 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	RU	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			
40	YU	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			

- Molecule 41 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	RV	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			
41	YV	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			

- Molecule 42 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	RW	113	Total	C	N	O	S	0	0	0
			900	566	177	155	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	YW	113	Total	C	N	O	S	0	0	0
			900	566	177	155	2			

- Molecule 43 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
43	RX	92	Total	C	N	O	0	0	0
			725	471	131	123			
43	YX	92	Total	C	N	O	0	0	0
			725	471	131	123			

- Molecule 44 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	RY	102	Total	C	N	O	S	0	0	0
			785	505	150	125	5			
44	YY	102	Total	C	N	O	S	0	0	0
			785	505	150	125	5			

- Molecule 45 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	RZ	183	Total	C	N	O	S	0	0	0
			1461	933	260	265	3			
45	YZ	183	Total	C	N	O	S	0	0	0
			1461	933	260	265	3			

- Molecule 46 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	R0	82	Total	C	N	O	S	0	0	0
			648	401	138	108	1			
46	Y0	82	Total	C	N	O	S	0	0	0
			648	401	138	108	1			

- Molecule 47 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	R1	97	Total	C	N	O	S	0	0	0
			763	481	150	131	1			
47	Y1	97	Total	C	N	O	S	0	0	0
			763	481	150	131	1			

- Molecule 48 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	R2	69	Total	C	N	O	S	0	0	0
			581	358	118	104	1			
48	Y2	69	Total	C	N	O	S	0	0	0
			581	358	118	104	1			

- Molecule 49 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
49	R3	59	Total	C	N	O	0	0	0
			469	298	90	81			
49	Y3	59	Total	C	N	O	0	0	0
			469	298	90	81			

- Molecule 50 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	R4	71	Total	C	N	O	S	0	0	0
			581	364	108	104	5			
50	Y4	71	Total	C	N	O	S	0	0	0
			581	364	108	104	5			

- Molecule 51 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	R5	59	Total	C	N	O	S	0	0	0
			459	288	90	76	5			
51	Y5	58	Total	C	N	O	S	0	0	0
			454	285	89	75	5			

- Molecule 52 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	R6	49	Total	C	N	O	S	0	0	0
			424	264	87	69	4			
52	Y6	49	Total	C	N	O	S	0	0	0
			424	264	87	69	4			

- Molecule 53 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	R7	49	Total	C	N	O	S	0	0	0
			430	263	108	57	2			
53	Y7	49	Total	C	N	O	S	0	0	0
			430	263	108	57	2			

- Molecule 54 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	R8	64	Total	C	N	O	S	0	0	0
			517	331	102	82	2			
54	Y8	64	Total	C	N	O	S	0	0	0
			517	331	102	82	2			

- Molecule 55 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	R9	37	Total	C	N	O	S	0	0	0
			307	188	68	47	4			
55	Y9	37	Total	C	N	O	S	0	0	0
			307	188	68	47	4			

- Molecule 56 is a RNA chain called CC-Puro.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
56	Z5	3	Total	C	N	O	P	0	0	0
			74	40	13	19	2			
56	Z6	3	Total	C	N	O	P	0	0	0
			74	40	13	19	2			

- Molecule 57 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	QA	95	Total	Mg	0	0
			95	95		
57	QE	1	Total	Mg	0	0
			1	1		
57	QF	1	Total	Mg	0	0
			1	1		
57	QM	1	Total	Mg	0	0
			1	1		
57	QV	3	Total	Mg	0	0
			3	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
57	QX	2	Total Mg 2 2	0	0
57	QY	1	Total Mg 1 1	0	0
57	RA	307	Total Mg 307 307	0	0
57	RB	5	Total Mg 5 5	0	0
57	RD	1	Total Mg 1 1	0	0
57	RE	2	Total Mg 2 2	0	0
57	RF	1	Total Mg 1 1	0	0
57	RR	1	Total Mg 1 1	0	0
57	R0	1	Total Mg 1 1	0	0
57	R1	1	Total Mg 1 1	0	0
57	R5	1	Total Mg 1 1	0	0
57	R8	1	Total Mg 1 1	0	0
57	XA	114	Total Mg 114 114	0	0
57	XF	1	Total Mg 1 1	0	0
57	XV	3	Total Mg 3 3	0	0
57	XX	2	Total Mg 2 2	0	0
57	YA	337	Total Mg 337 337	0	0
57	YB	3	Total Mg 3 3	0	0
57	YD	1	Total Mg 1 1	0	0
57	YE	2	Total Mg 2 2	0	0
57	YP	2	Total Mg 2 2	0	0

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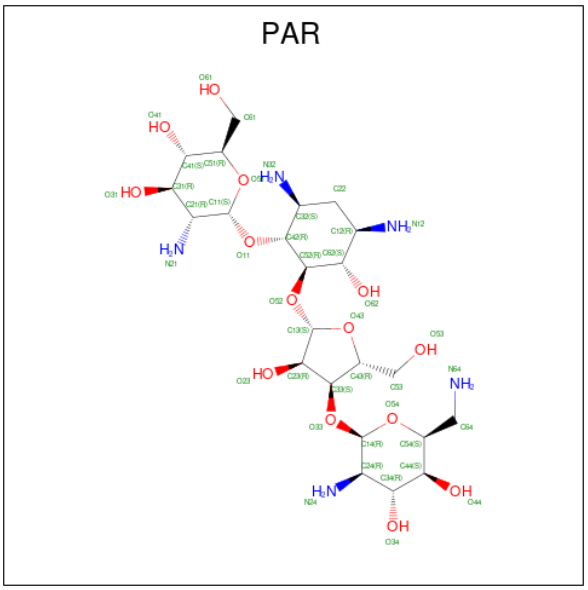
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
57	YQ	1	Total Mg 1 1	0	0
57	Y0	3	Total Mg 3 3	0	0
57	Y5	1	Total Mg 1 1	0	0
57	Y7	1	Total Mg 1 1	0	0

- Molecule 58 is PAROMOMYCIN (CCD ID: PAR) (formula: $C_{23}H_{45}N_5O_{14}$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
58	QA	1	Total C N O 42 23 5 14	0	0
58	XA	1	Total C N O 42 23 5 14	0	0

- Molecule 59 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
59	QD	1	Total Zn 1 1	0	0
59	QN	1	Total Zn 1 1	0	0
59	XD	1	Total Zn 1 1	0	0

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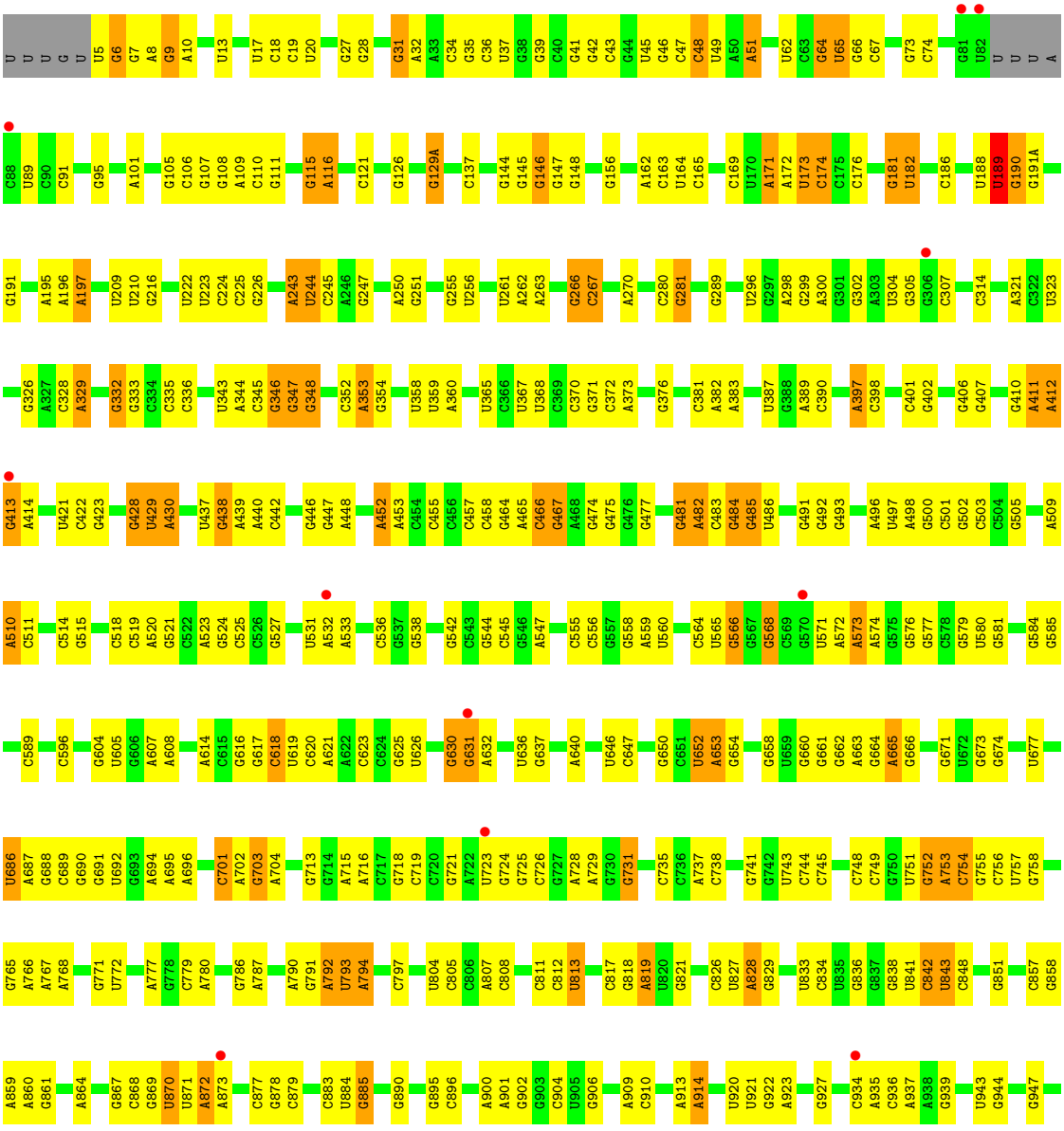
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
59	XN	1	1	1	0	0

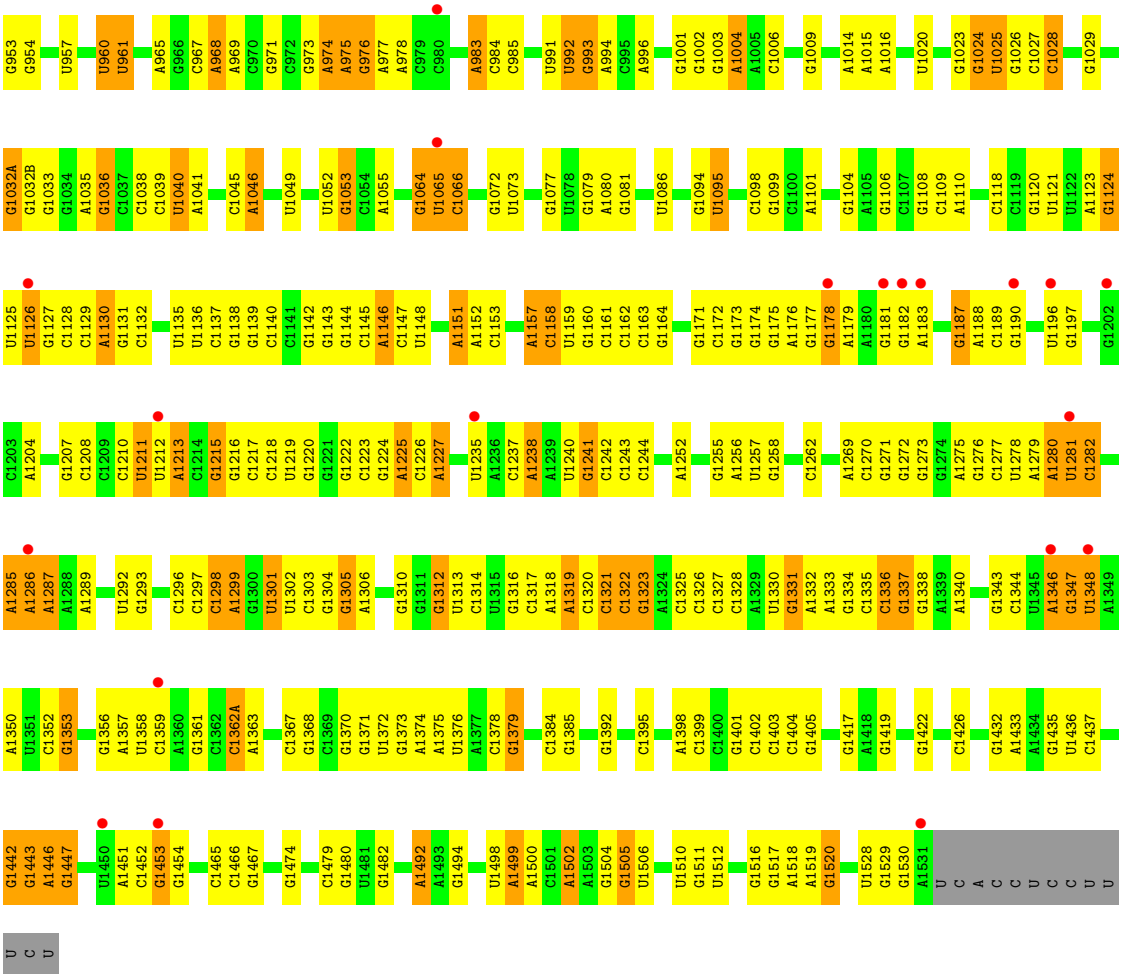
3 Residue-property plots (i)

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

- Molecule 1: 16S rRNA

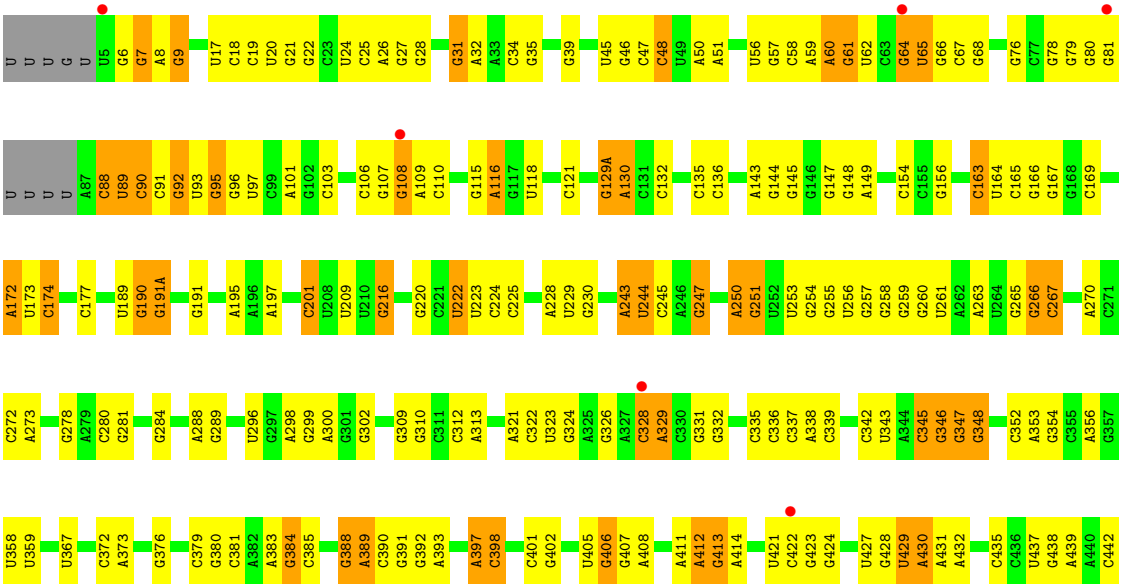
Chain QA:

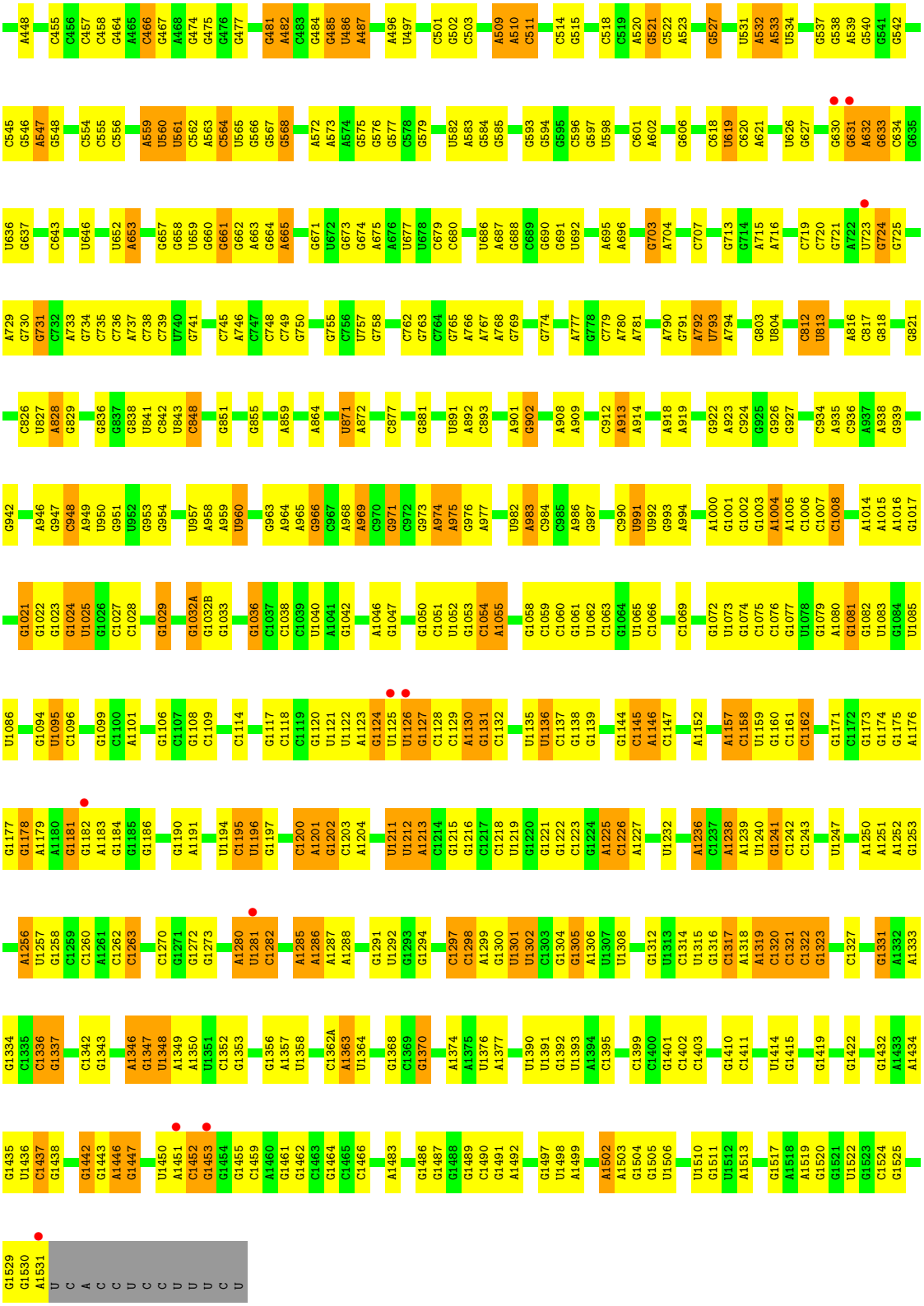




• Molecule 1: 16S rRNA

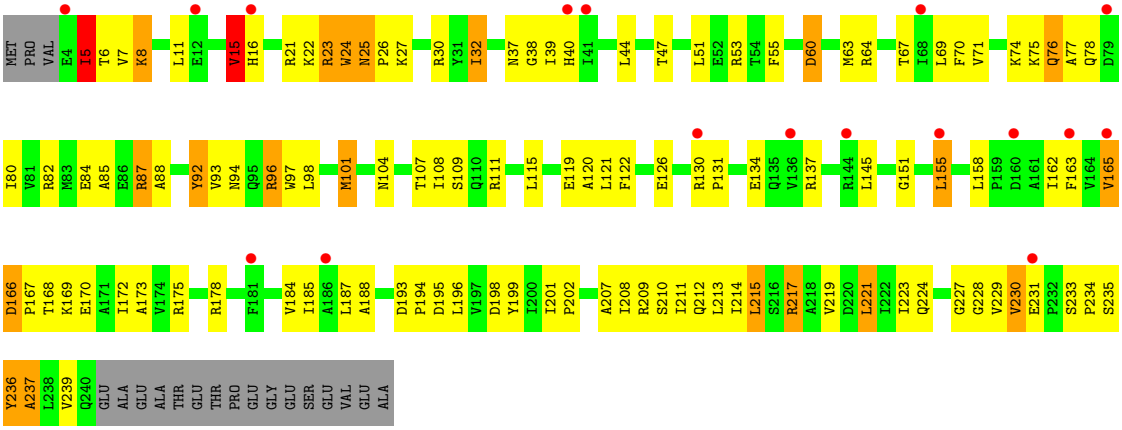
Chain XA:





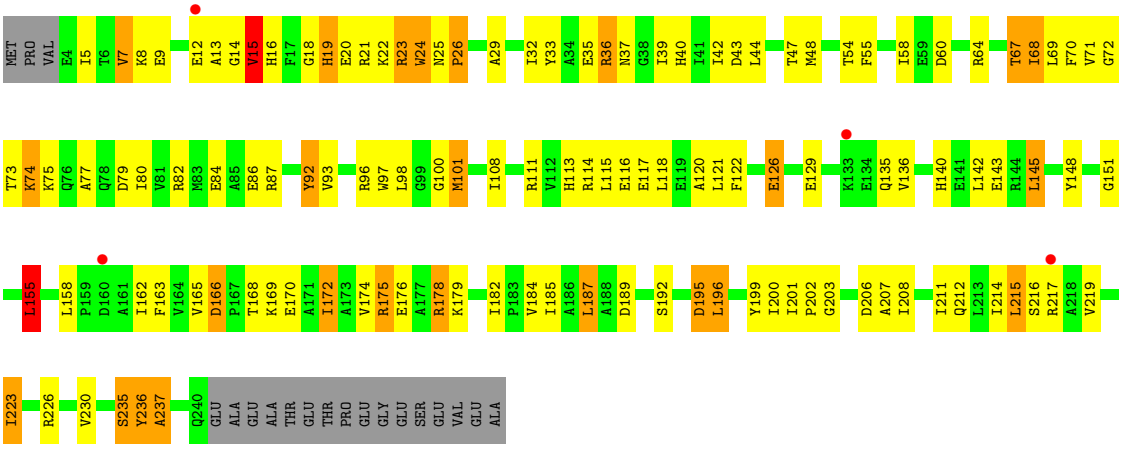
• Molecule 2: 30S ribosomal protein S2

Chain QB:



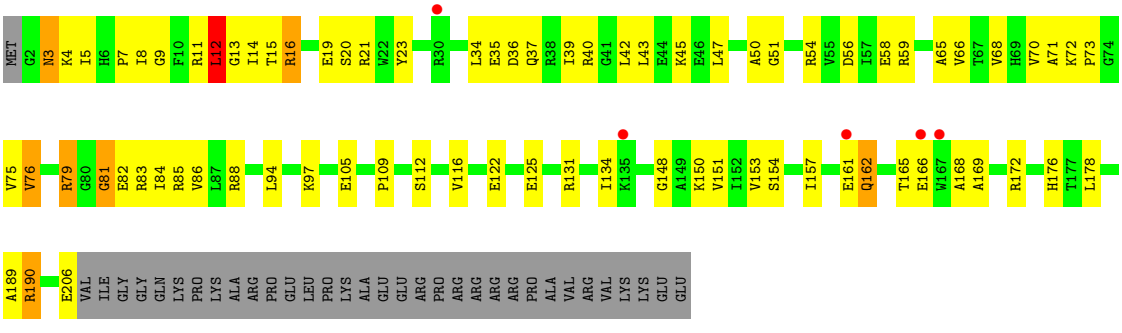
• Molecule 2: 30S ribosomal protein S2

Chain XB:



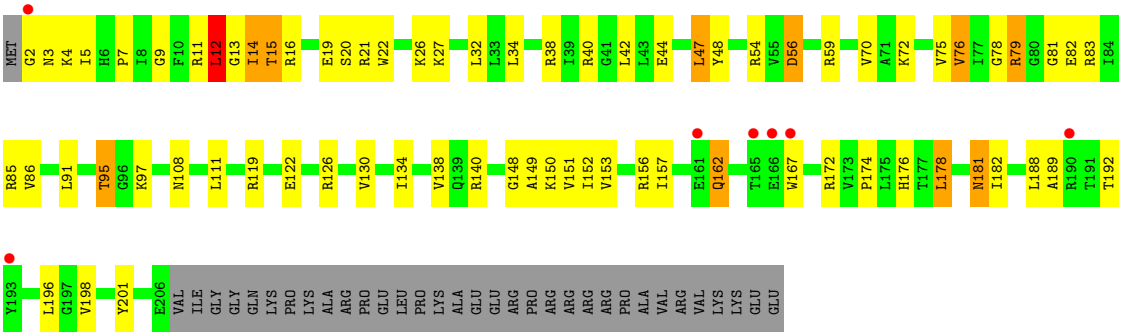
• Molecule 3: 30S ribosomal protein S3

Chain QC:

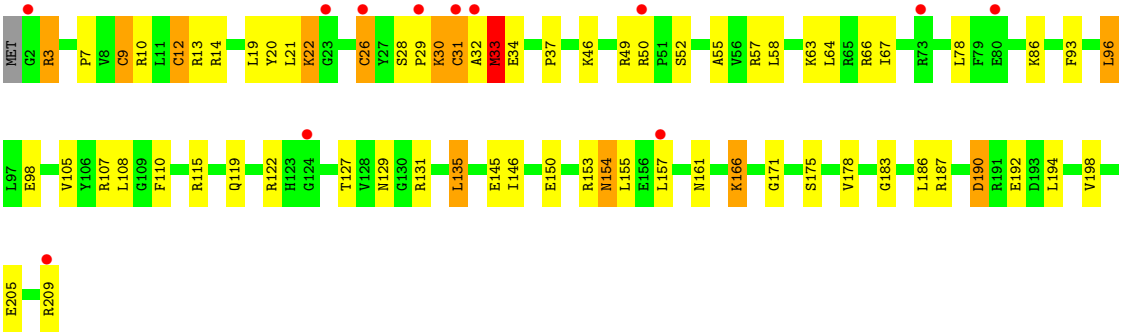


• Molecule 3: 30S ribosomal protein S3

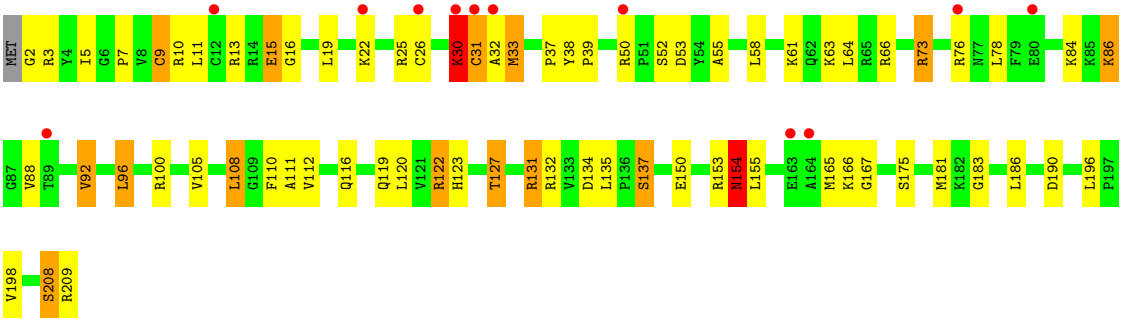
Chain XC:



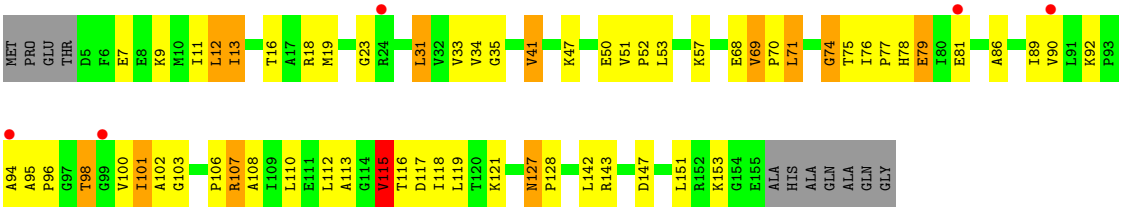
• Molecule 4: 30S ribosomal protein S4



• Molecule 4: 30S ribosomal protein S4

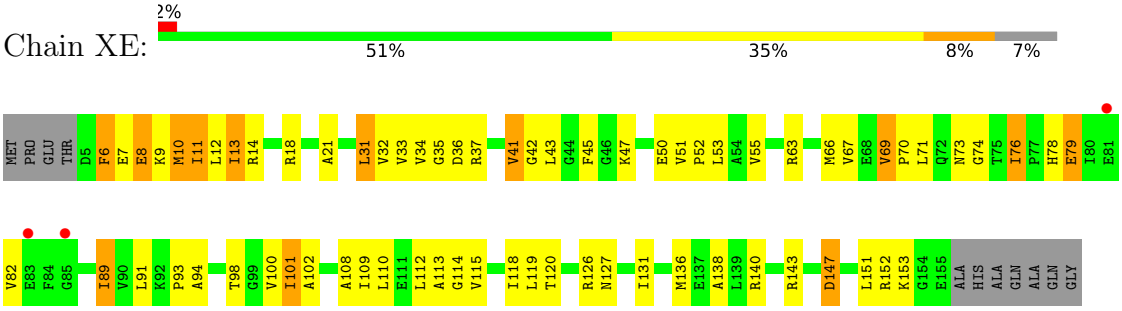


• Molecule 5: 30S ribosomal protein S5



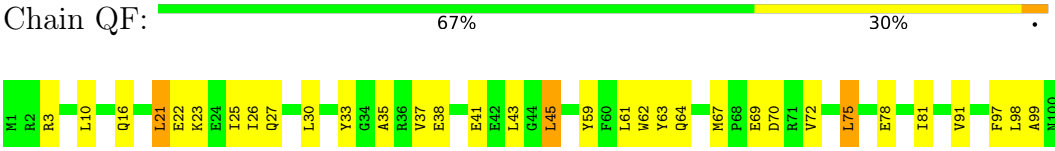
• Molecule 5: 30S ribosomal protein S5

Chain XE:



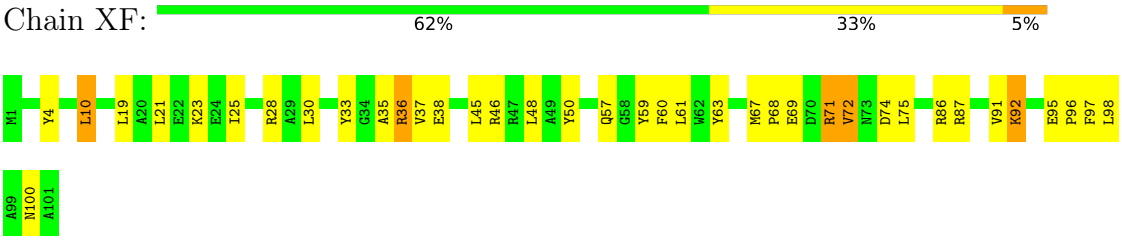
• Molecule 6: 30S ribosomal protein S6

Chain QF:



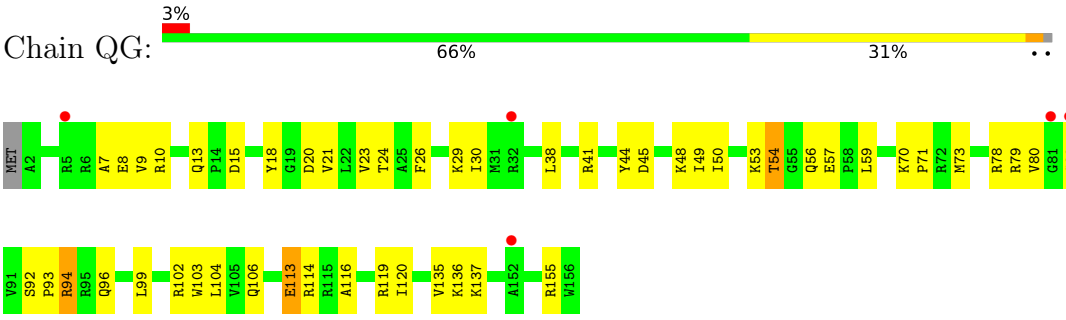
• Molecule 6: 30S ribosomal protein S6

Chain XF:



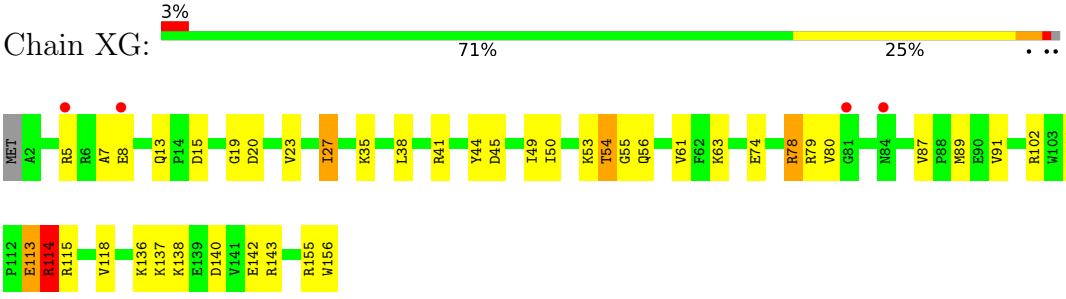
• Molecule 7: 30S ribosomal protein S7

Chain QG:



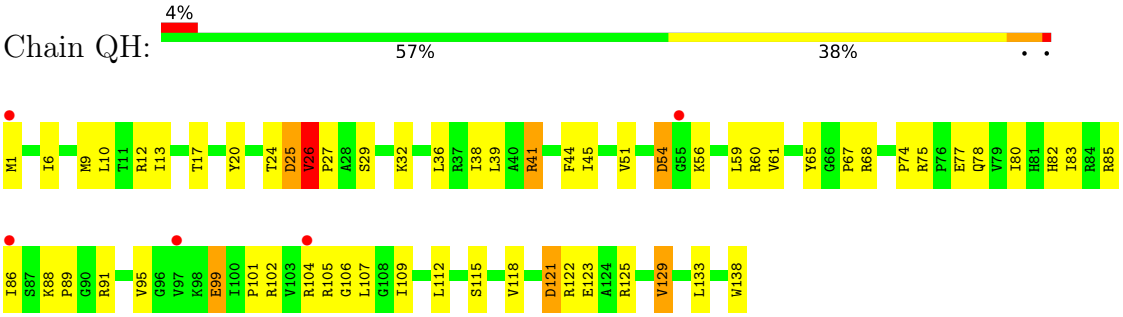
• Molecule 7: 30S ribosomal protein S7

Chain XG:



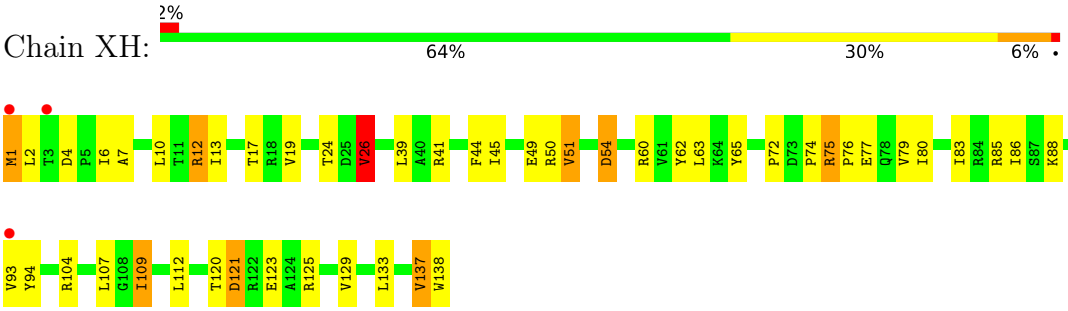
• Molecule 8: 30S ribosomal protein S8

Chain QH:



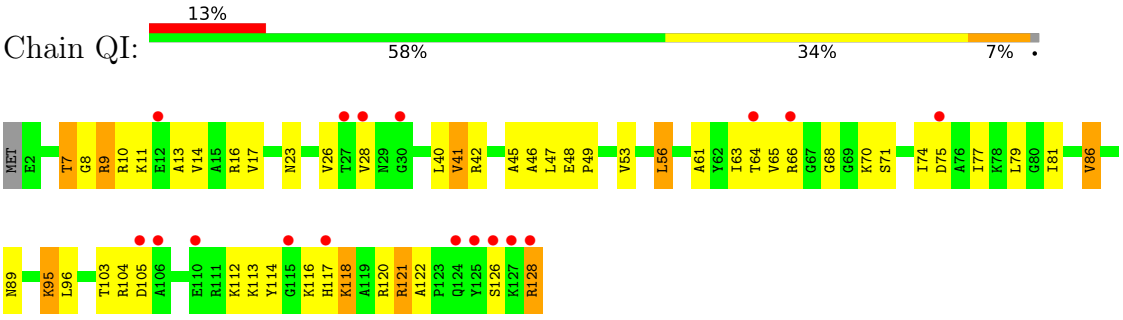
• Molecule 8: 30S ribosomal protein S8

Chain XH:



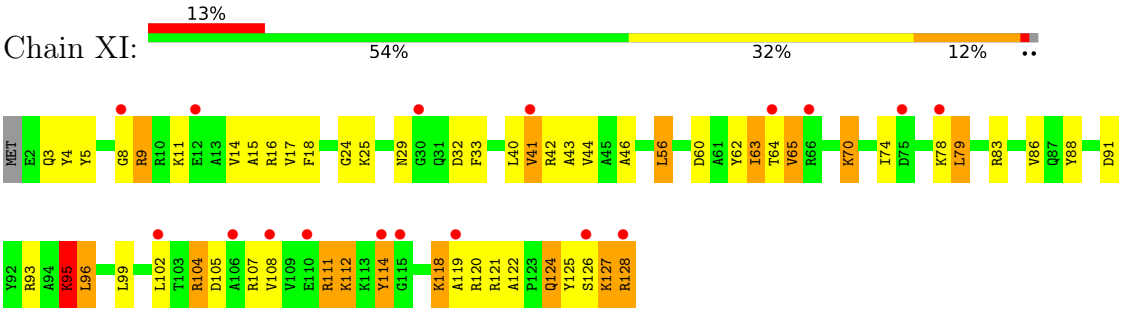
• Molecule 9: 30S ribosomal protein S9

Chain QI:



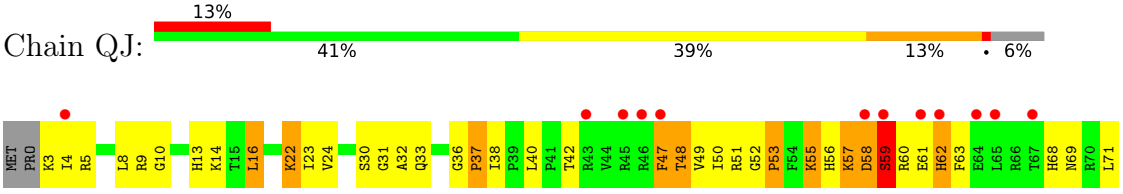
• Molecule 9: 30S ribosomal protein S9

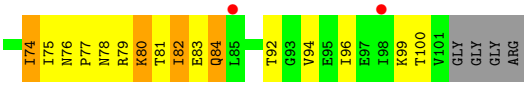
Chain XI:



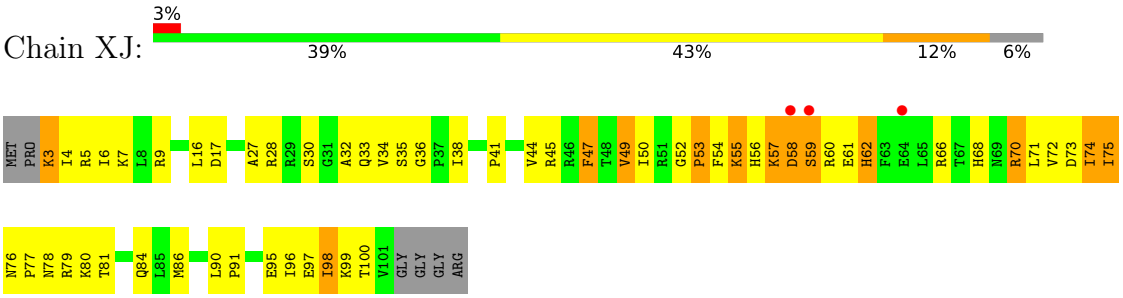
• Molecule 10: 30S ribosomal protein S10

Chain QJ:

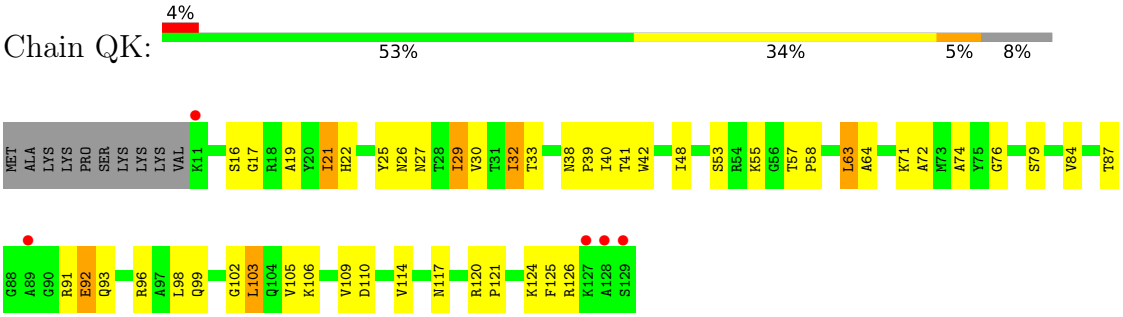




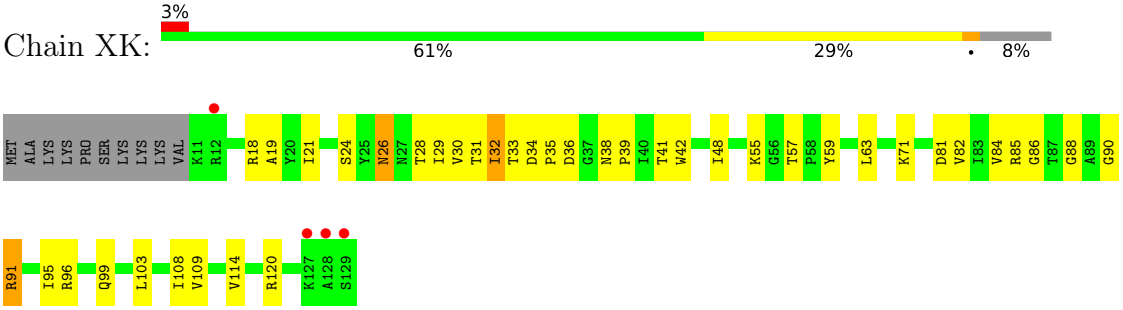
• Molecule 10: 30S ribosomal protein S10



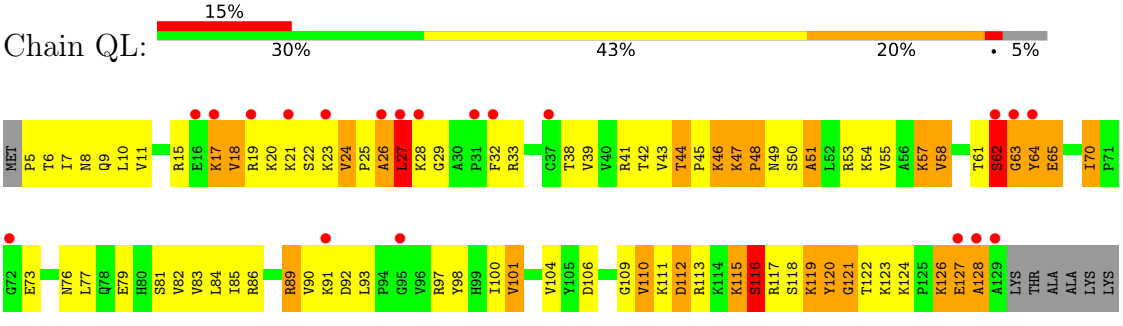
• Molecule 11: 30S ribosomal protein S11



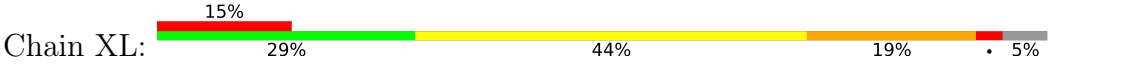
• Molecule 11: 30S ribosomal protein S11

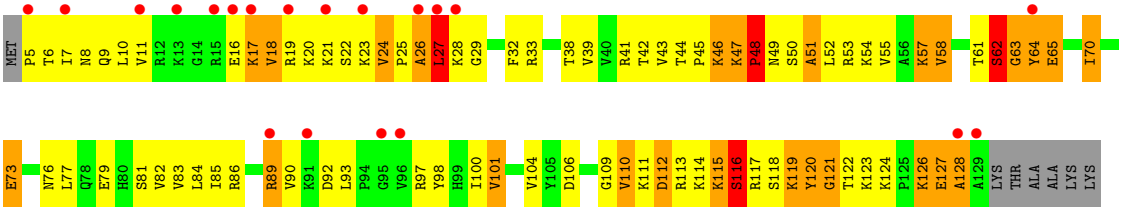


• Molecule 12: 30S ribosomal protein S12

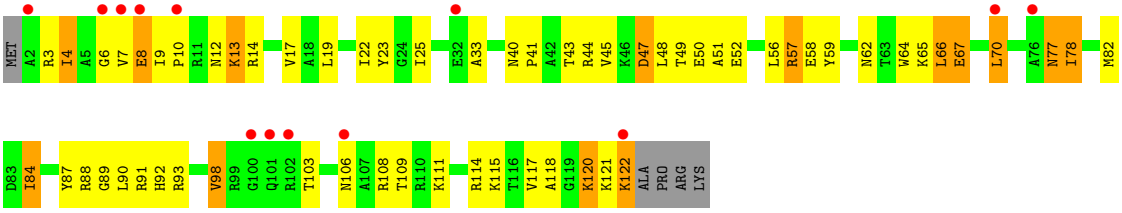


• Molecule 12: 30S ribosomal protein S12

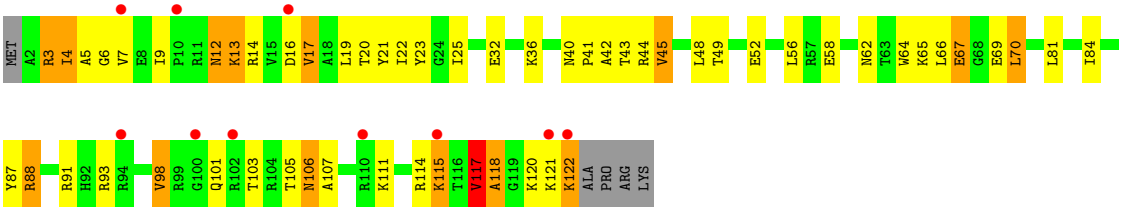




• Molecule 13: 30S ribosomal protein S13

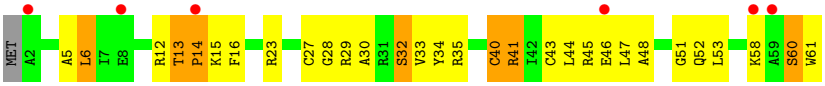


• Molecule 13: 30S ribosomal protein S13

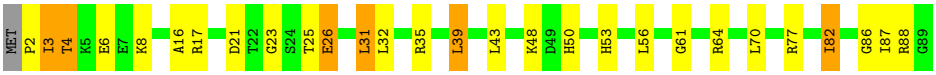


• Molecule 14: 30S ribosomal protein S14 type Z

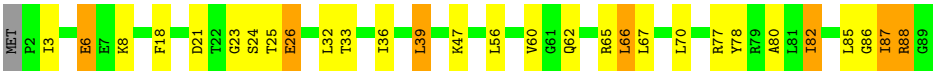
• Molecule 14: 30S ribosomal protein S14 type Z



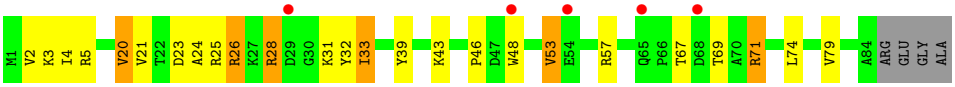
• Molecule 15: 30S ribosomal protein S15



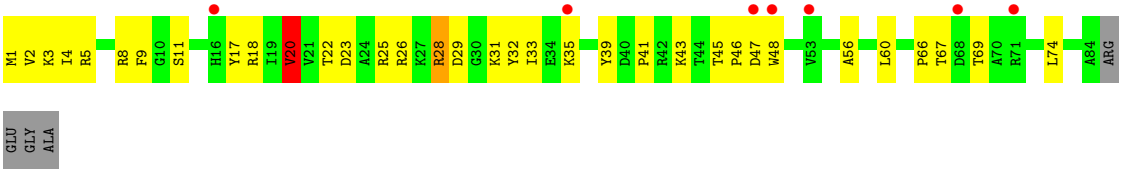
• Molecule 15: 30S ribosomal protein S15



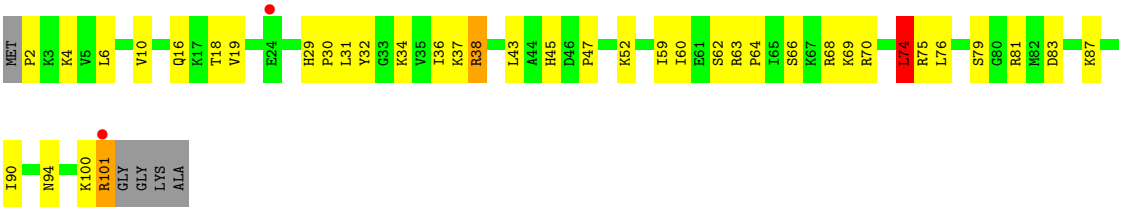
• Molecule 16: 30S ribosomal protein S16



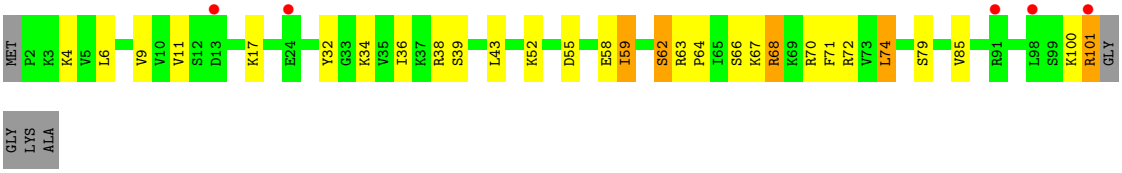
• Molecule 16: 30S ribosomal protein S16



• Molecule 17: 30S ribosomal protein S17

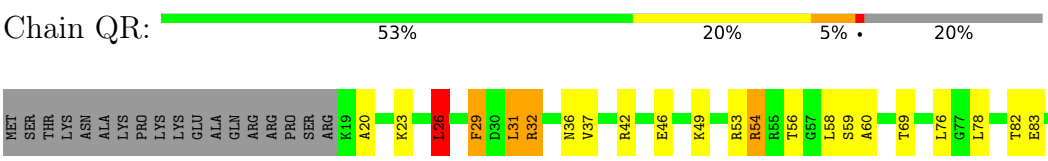


• Molecule 17: 30S ribosomal protein S17



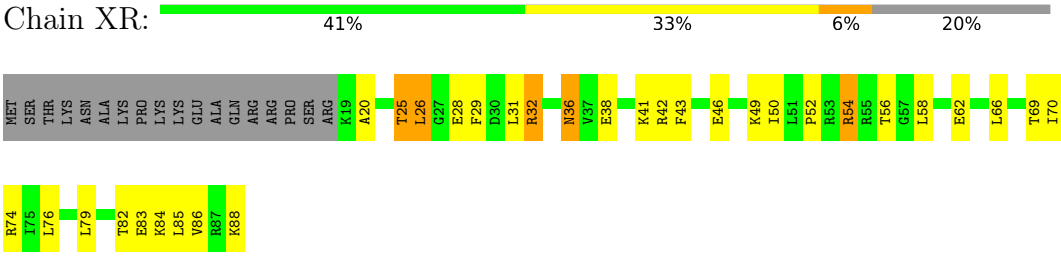
• Molecule 18: 30S ribosomal protein S18

Chain QR:



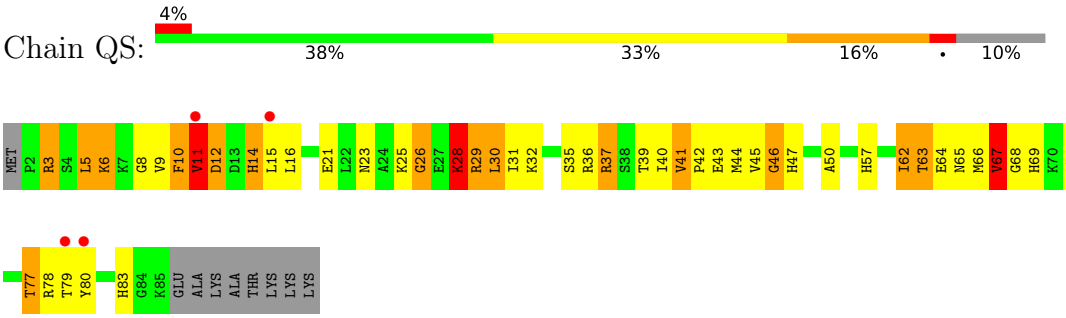
• Molecule 18: 30S ribosomal protein S18

Chain XR:



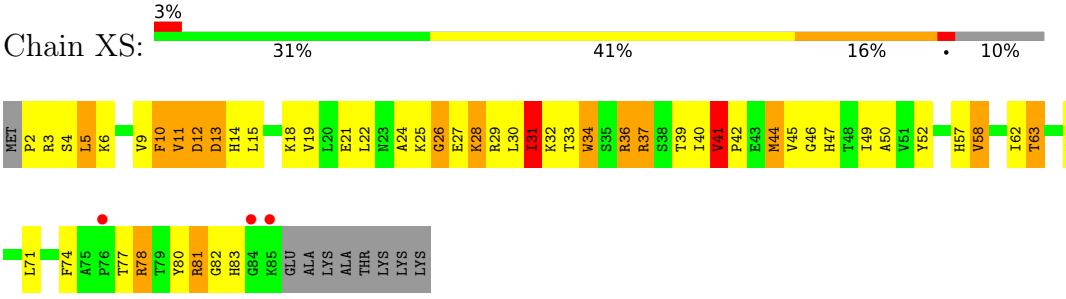
• Molecule 19: 30S ribosomal protein S19

Chain QS:



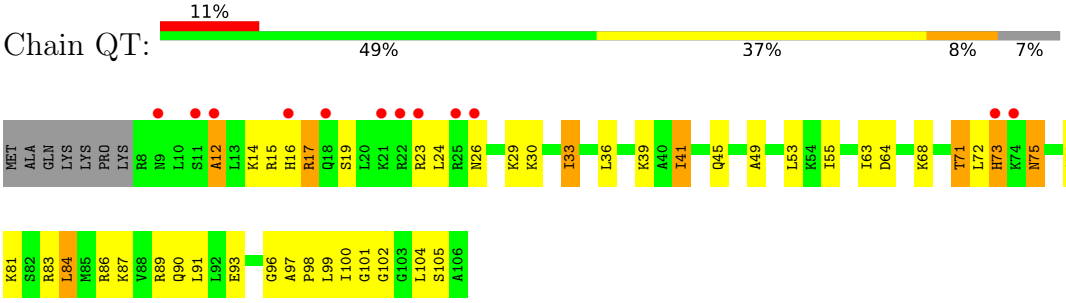
• Molecule 19: 30S ribosomal protein S19

Chain XS:



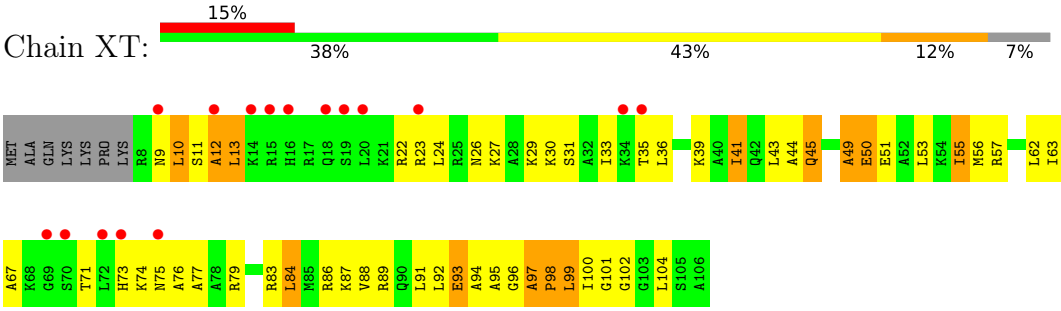
• Molecule 20: 30S ribosomal protein S20

Chain QT:



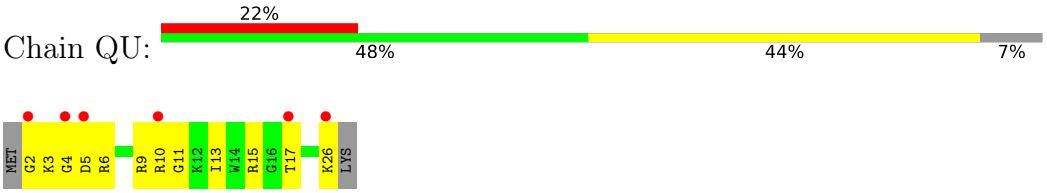
• Molecule 20: 30S ribosomal protein S20

Chain XT:



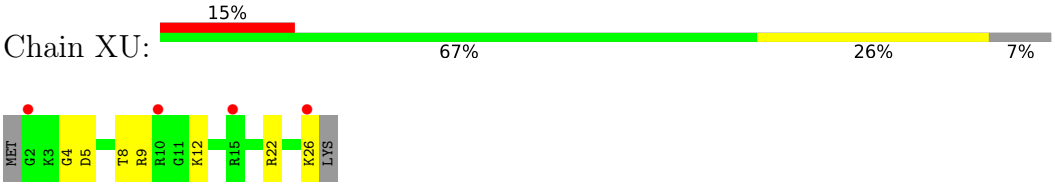
• Molecule 21: 30S ribosomal protein Thx

Chain QU:



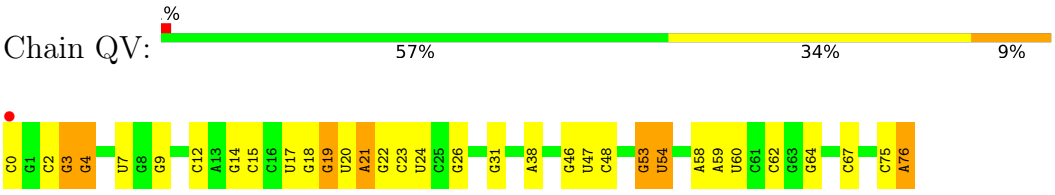
• Molecule 21: 30S ribosomal protein Thx

Chain XU:



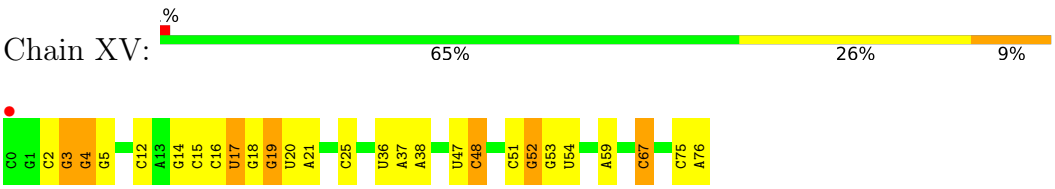
• Molecule 22: P-site tRNA-fMet

Chain QV:



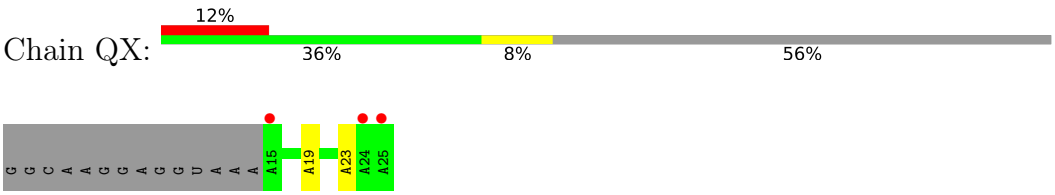
• Molecule 22: P-site tRNA-fMet

Chain XV:



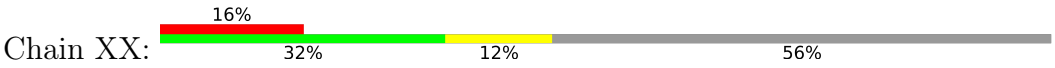
• Molecule 23: messenger RNA

Chain QX:



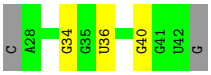
• Molecule 23: messenger RNA

Chain XX:



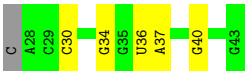
• Molecule 24: A-site ASL-Thr

Chain QY:



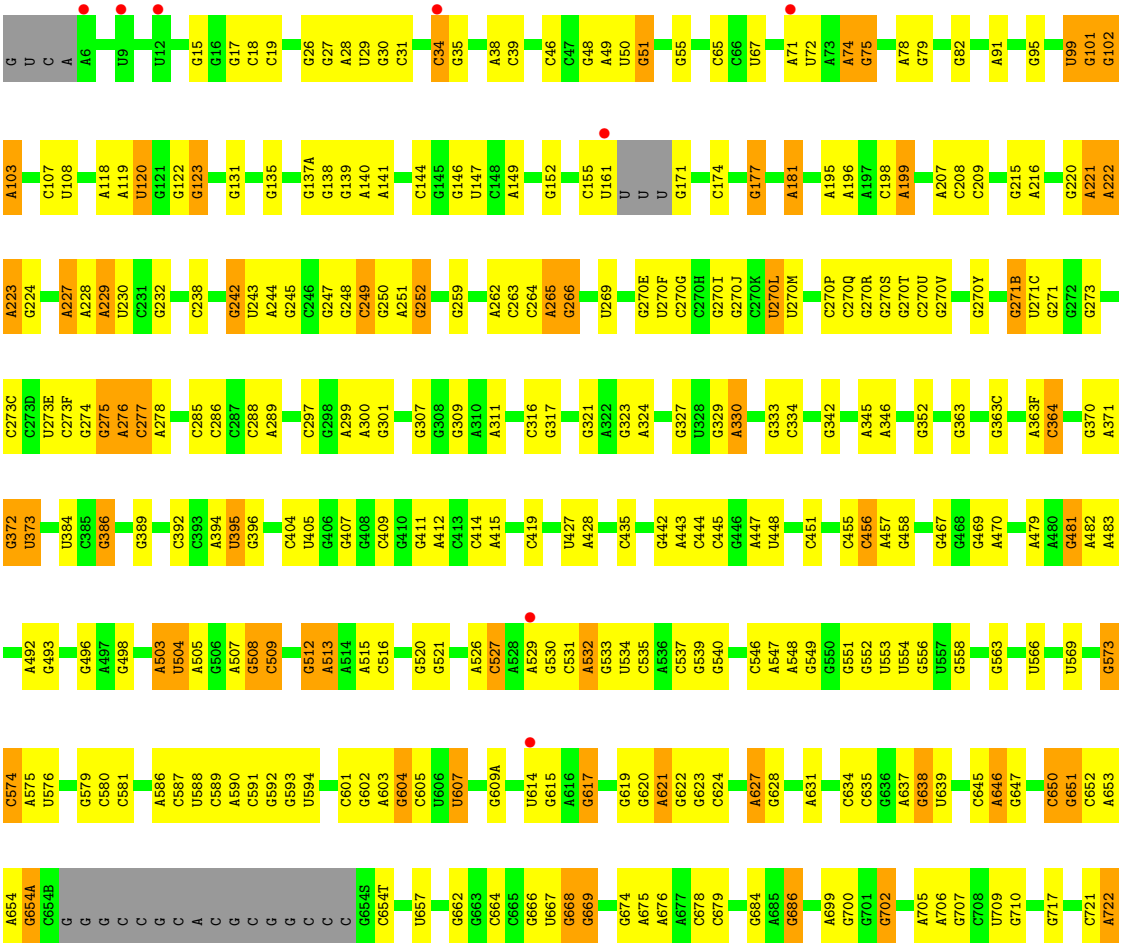
• Molecule 24: A-site ASL-Thr

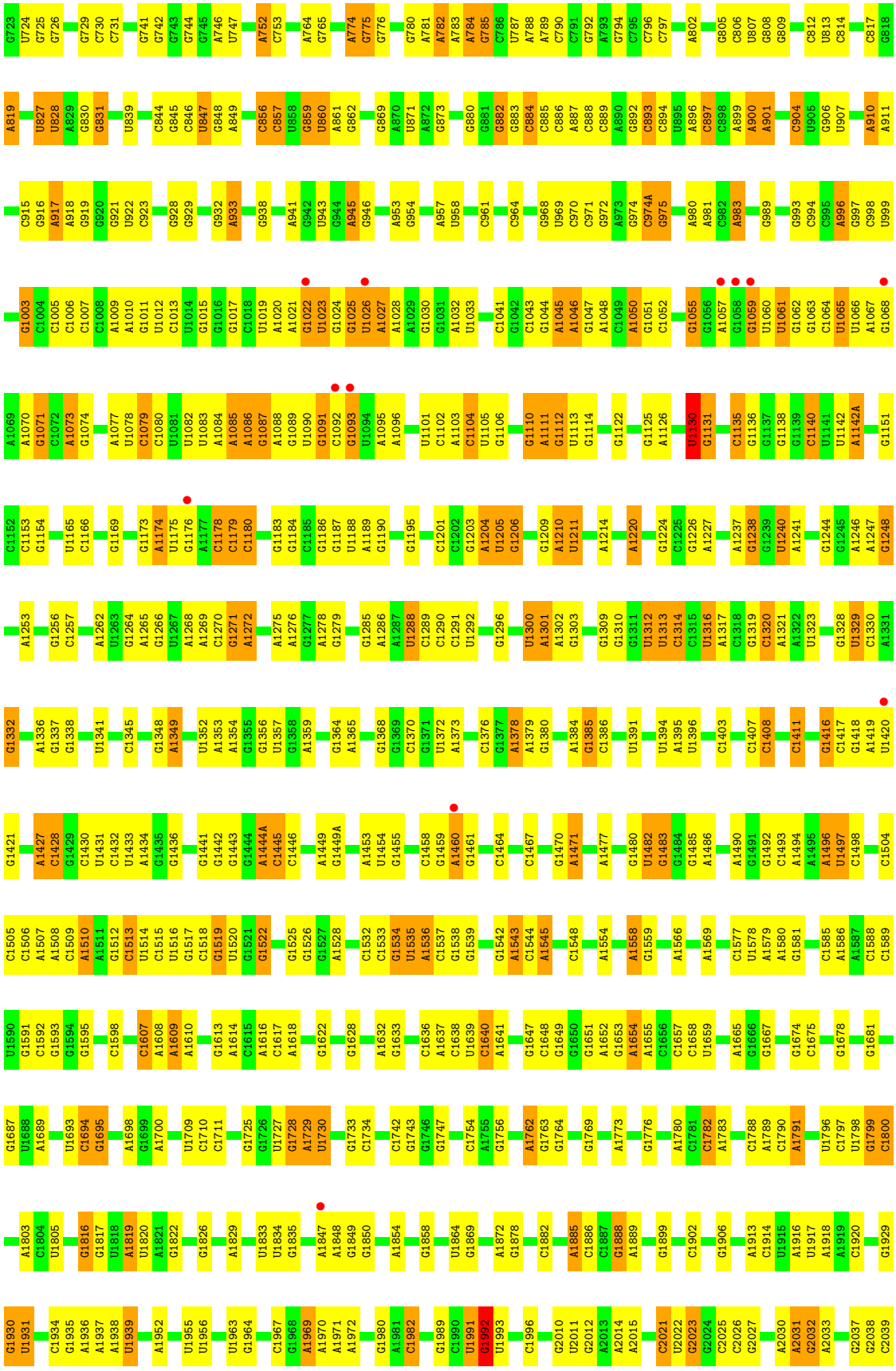
Chain XY:



• Molecule 25: 23S rRNA

Chain RA:

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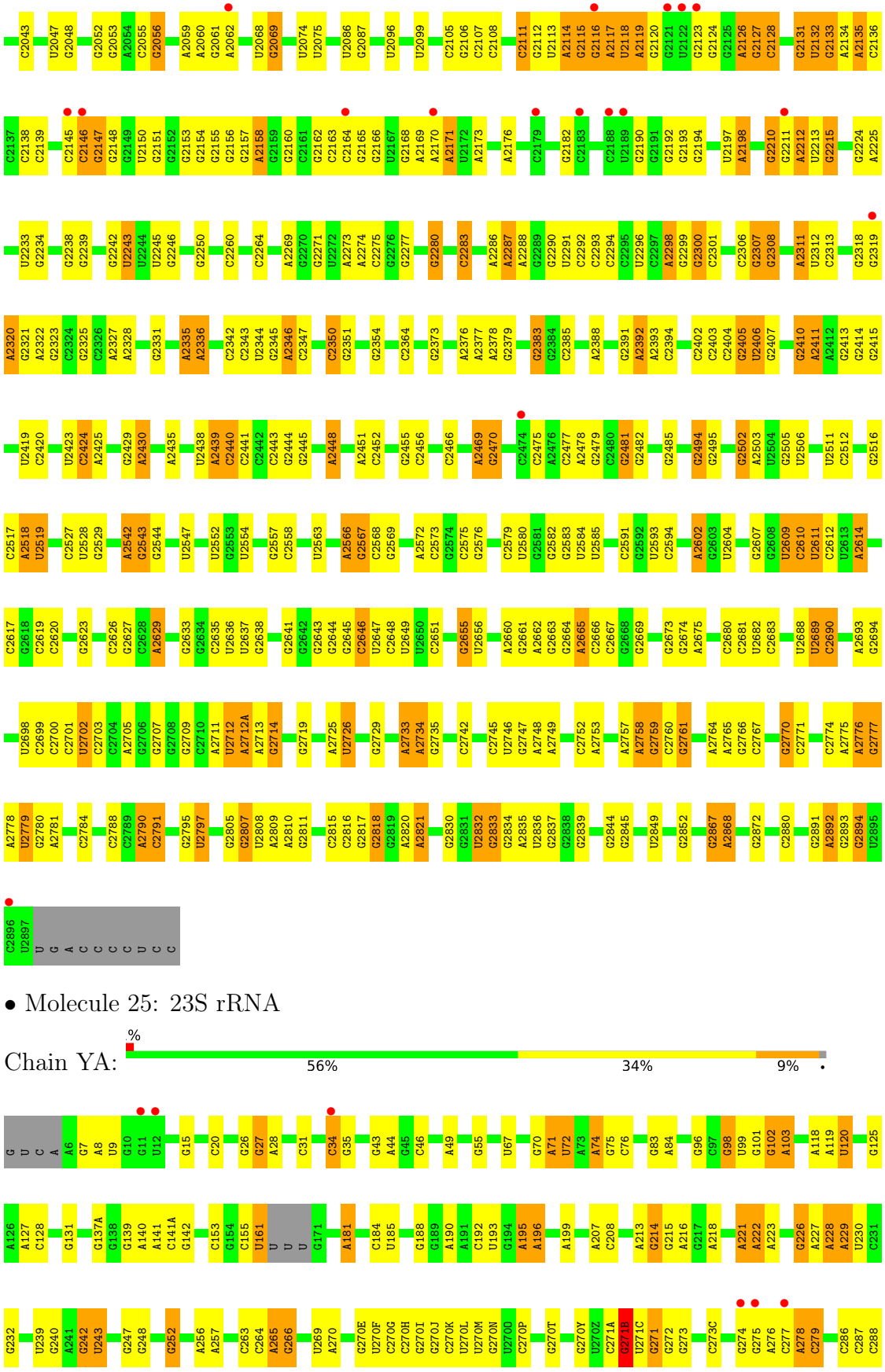
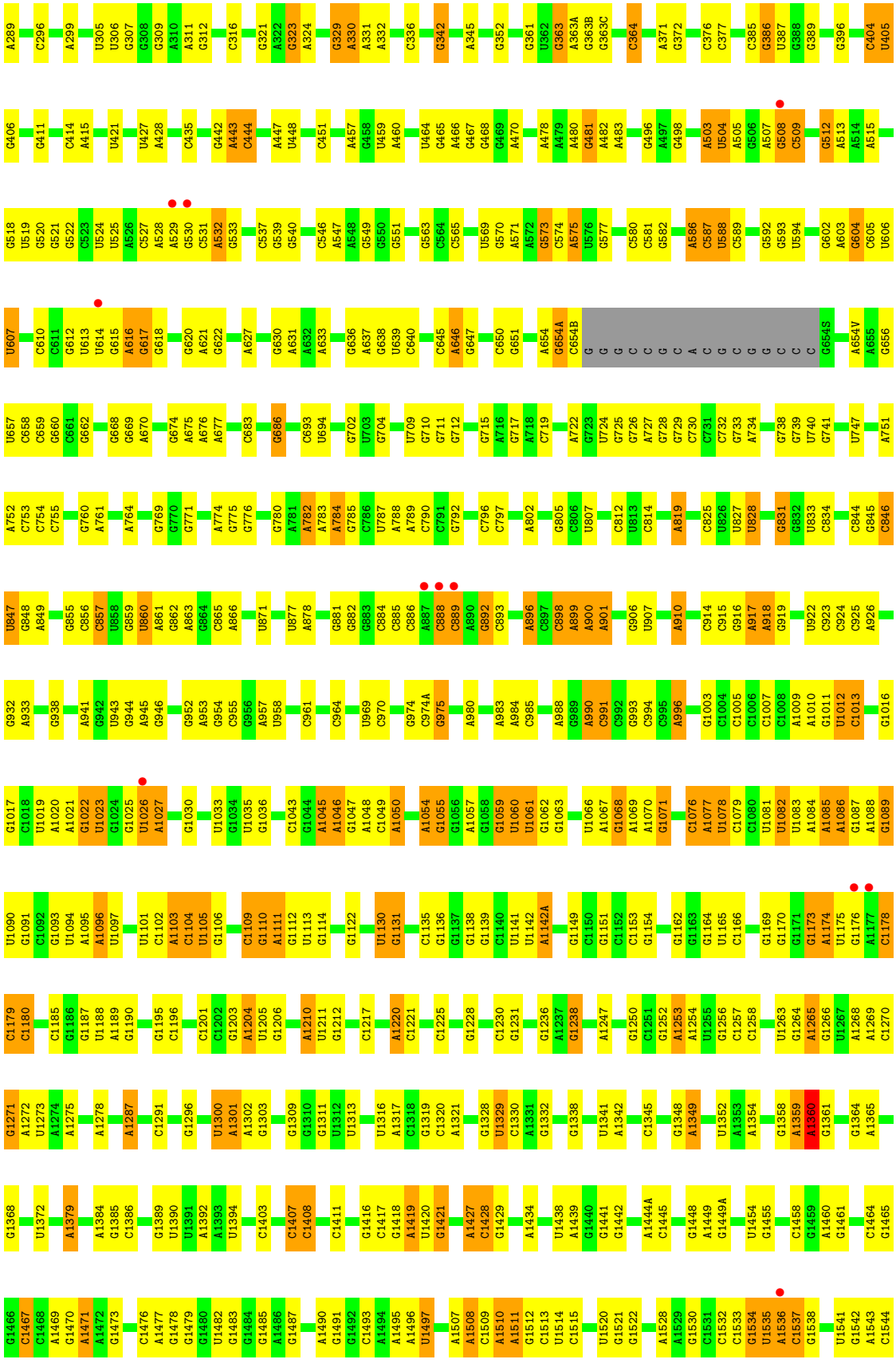


• Molecule 25: 23S rRNA

Chain YA: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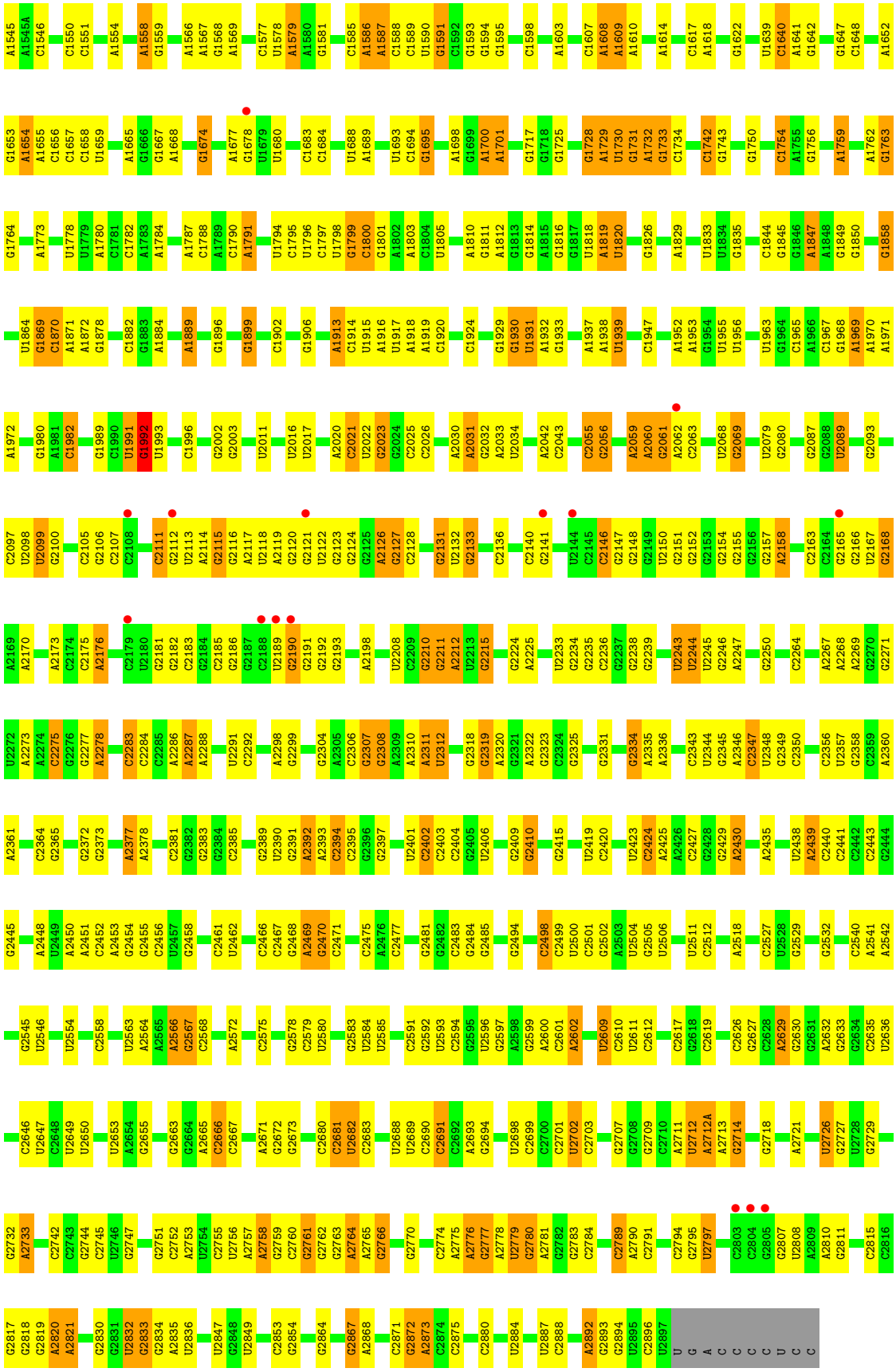
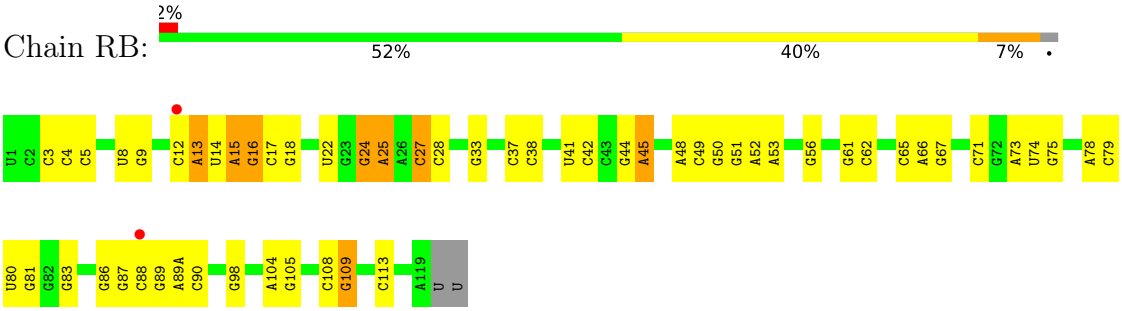


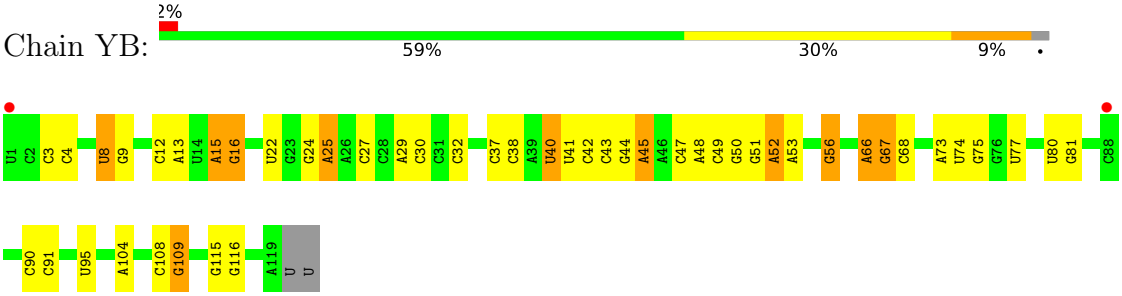
• Molecule 26: 5S rRNA

Chain RB:



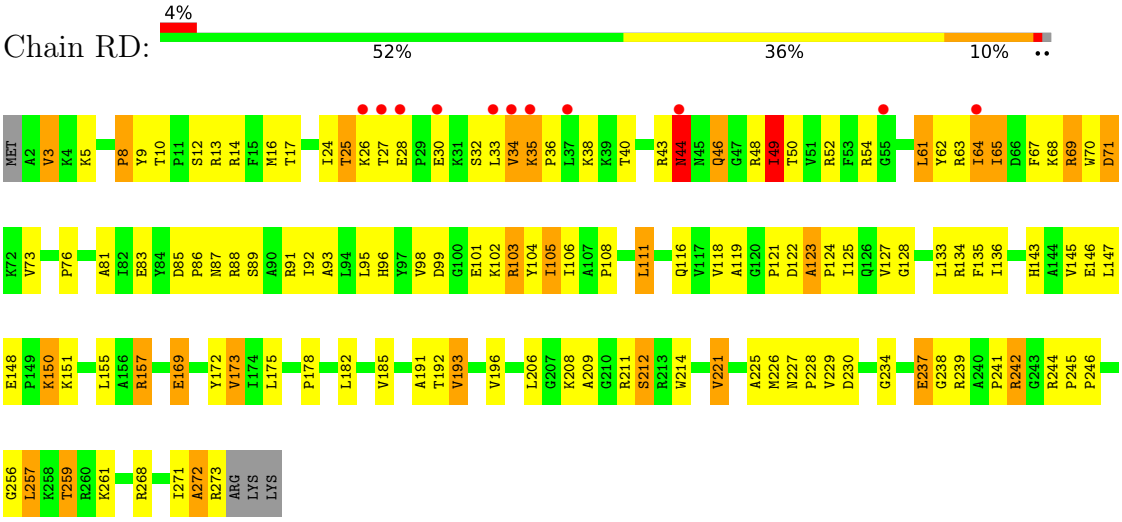
• Molecule 26: 5S rRNA

Chain YB:



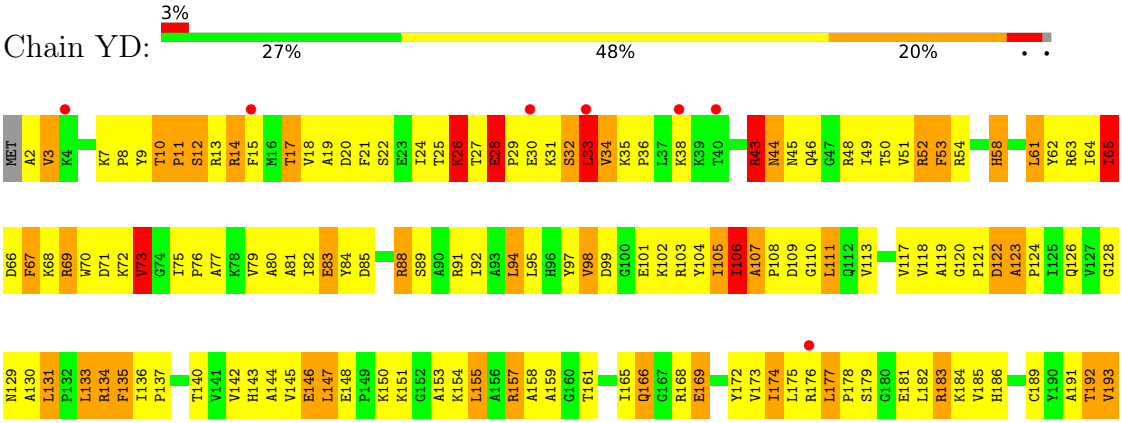
• Molecule 27: 50S ribosomal protein L2

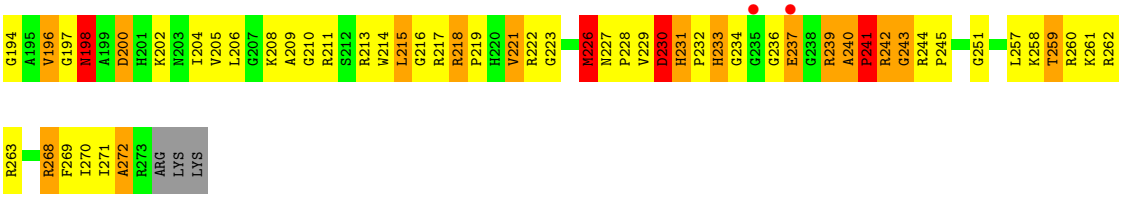
Chain RD:



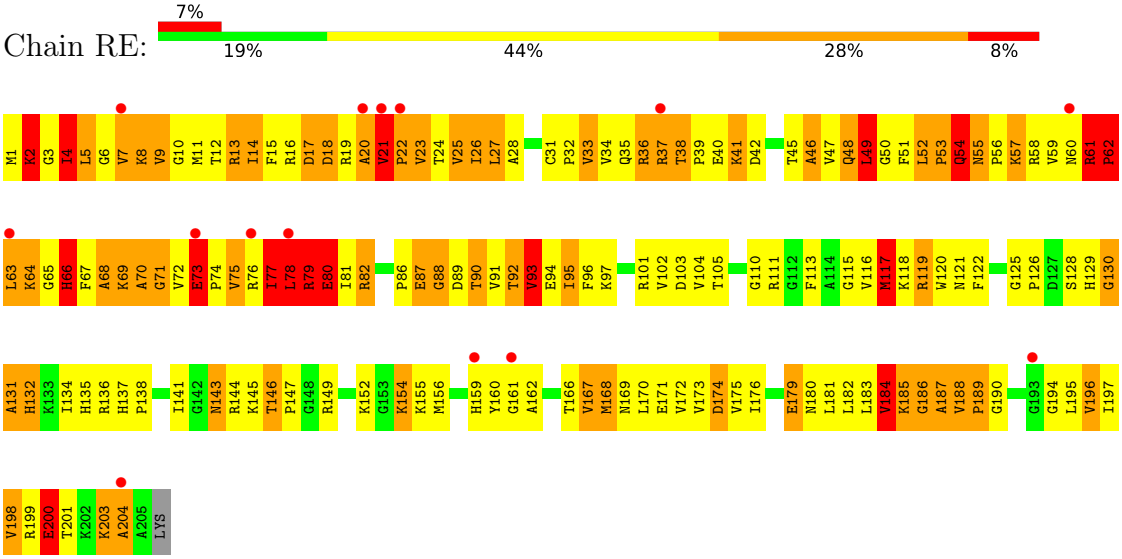
• Molecule 27: 50S ribosomal protein L2

Chain YD:

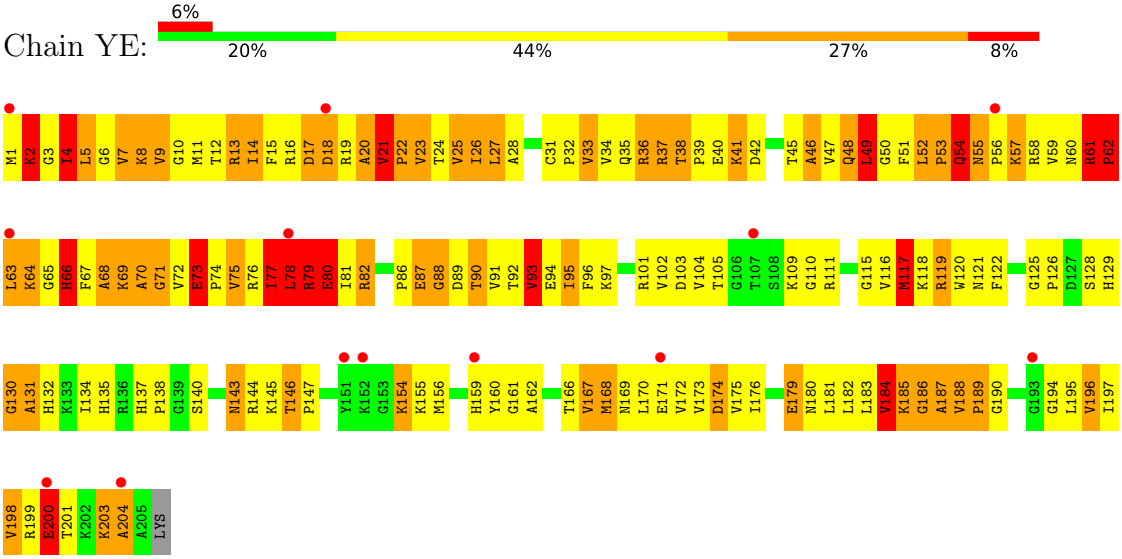




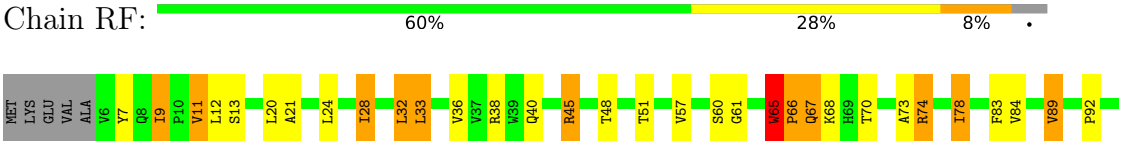
• Molecule 28: 50S ribosomal protein L3

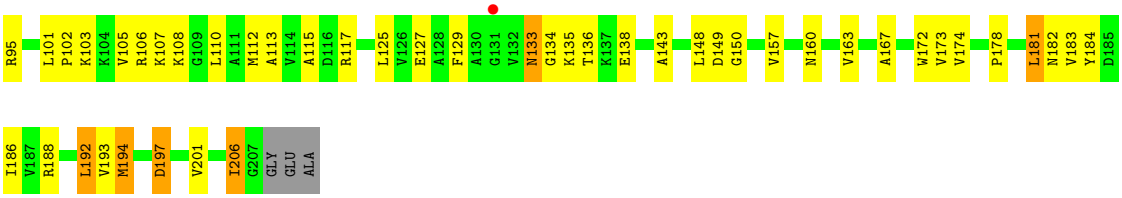


• Molecule 28: 50S ribosomal protein L3



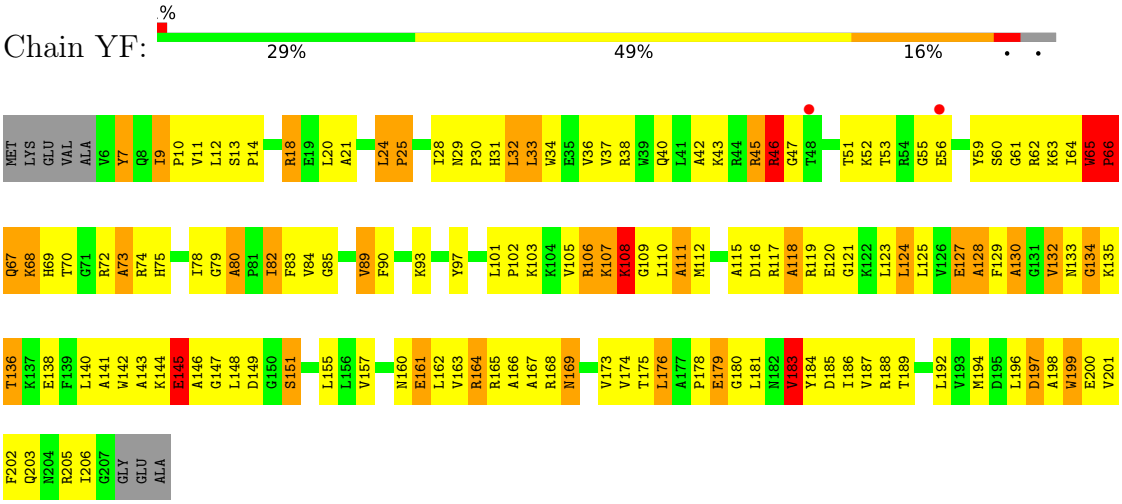
• Molecule 29: 50S ribosomal protein L4





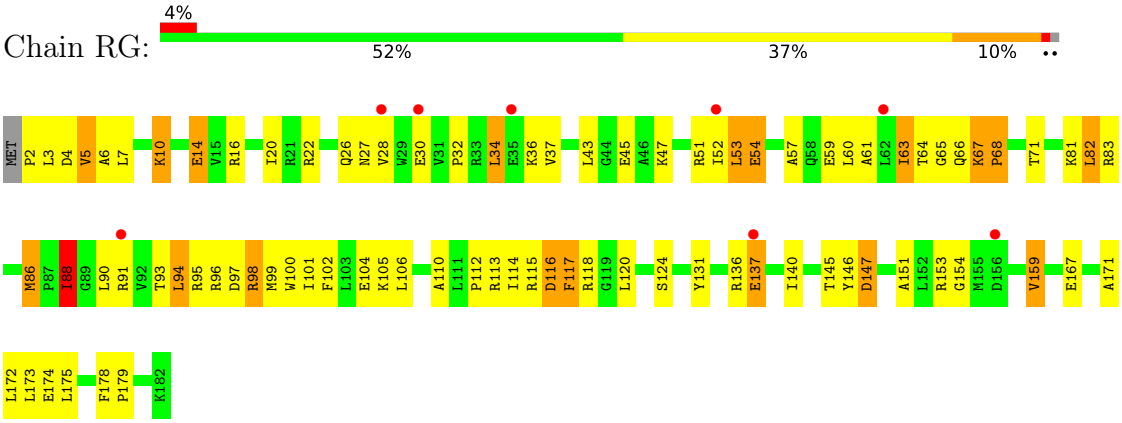
• Molecule 29: 50S ribosomal protein L4

Chain YF:



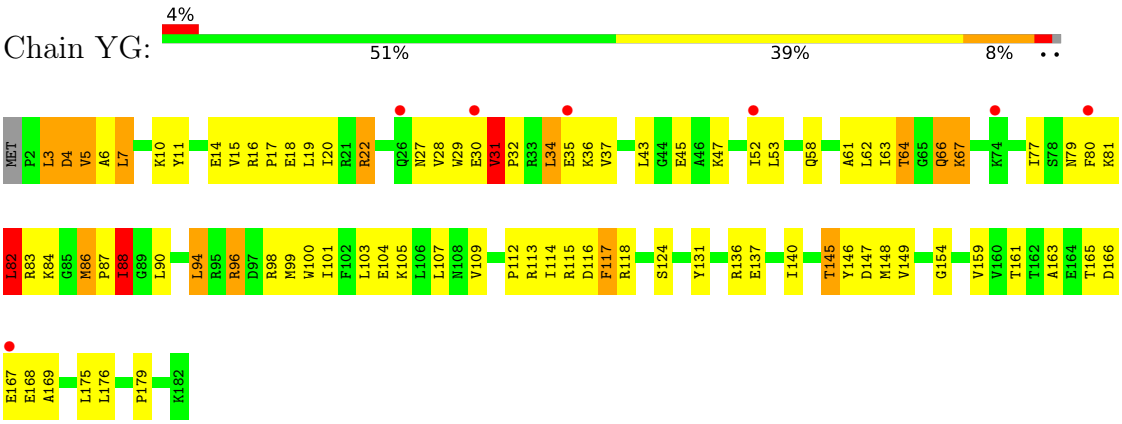
• Molecule 30: 50S ribosomal protein L5

Chain RG:



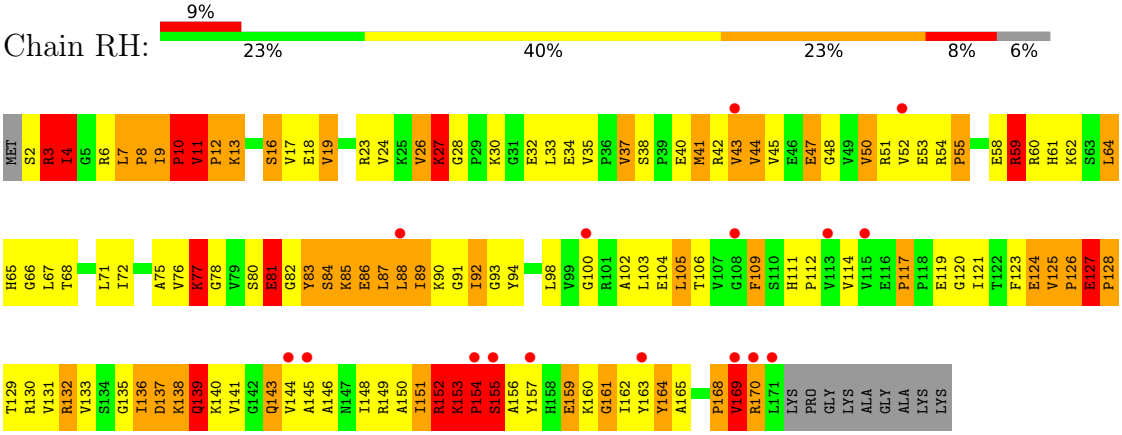
• Molecule 30: 50S ribosomal protein L5

Chain YG:



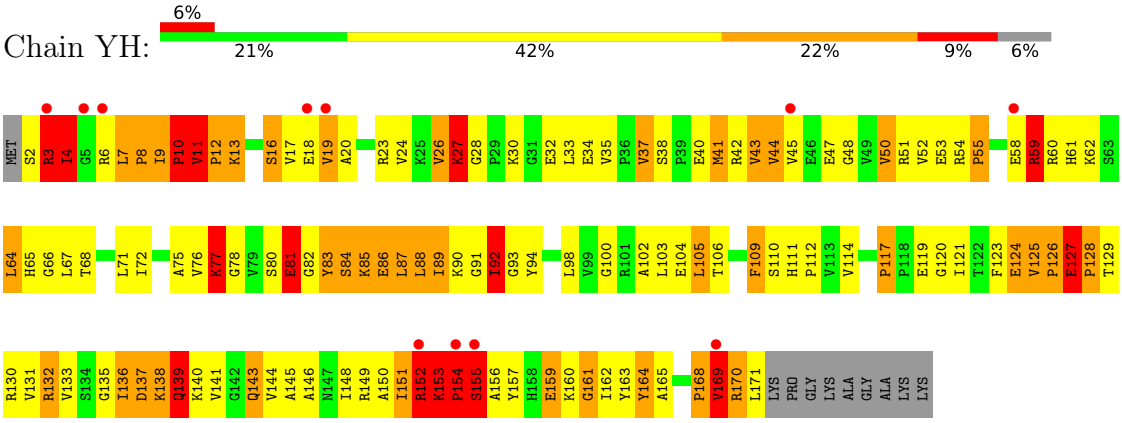
• Molecule 31: 50S ribosomal protein L6

Chain RH:



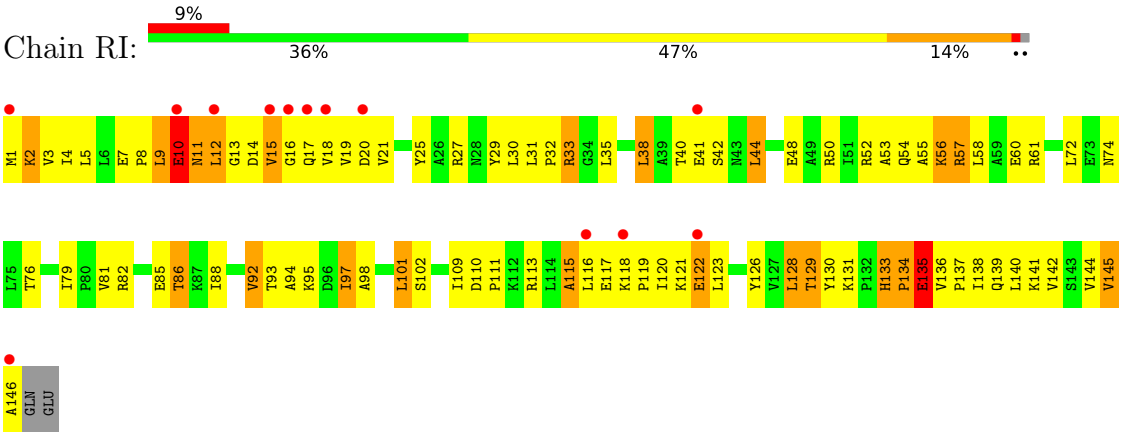
• Molecule 31: 50S ribosomal protein L6

Chain YH:



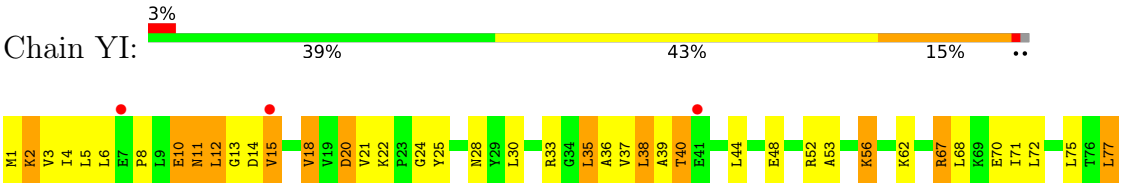
• Molecule 32: 50S ribosomal protein L9

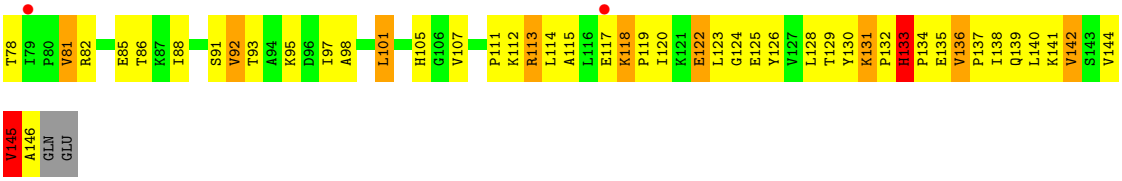
Chain RI:



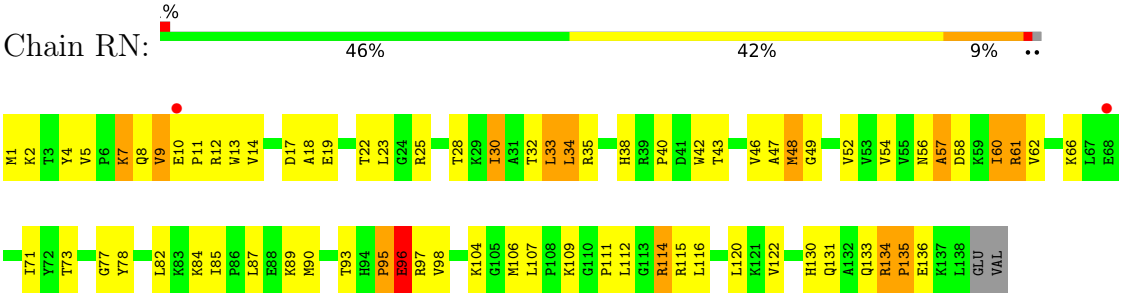
• Molecule 32: 50S ribosomal protein L9

Chain YI:

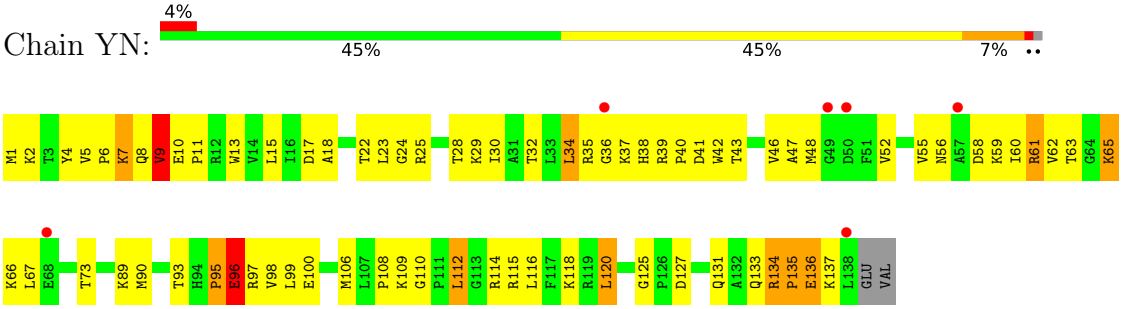




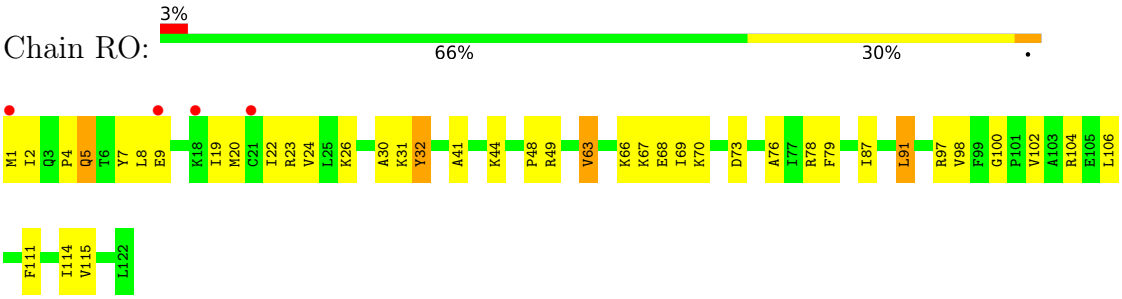
• Molecule 33: 50S ribosomal protein L13



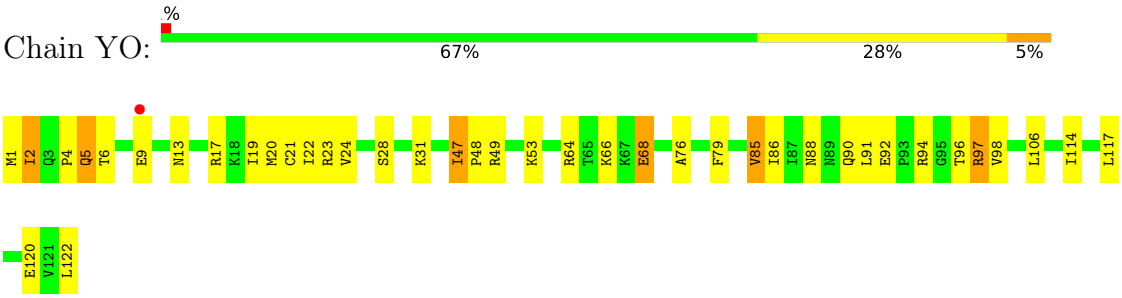
• Molecule 33: 50S ribosomal protein L13



• Molecule 34: 50S ribosomal protein L14

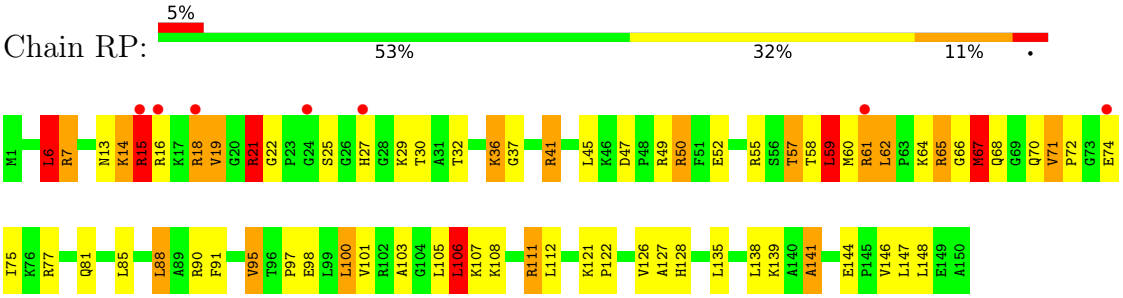


• Molecule 34: 50S ribosomal protein L14



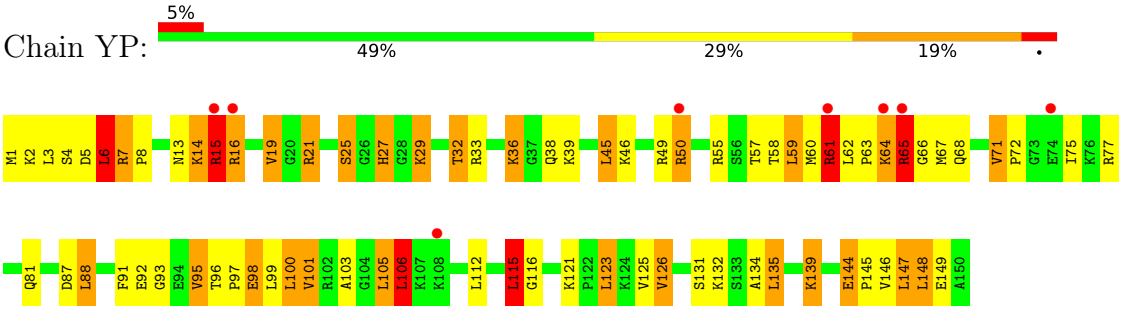
• Molecule 35: 50S ribosomal protein L15

Chain RP:



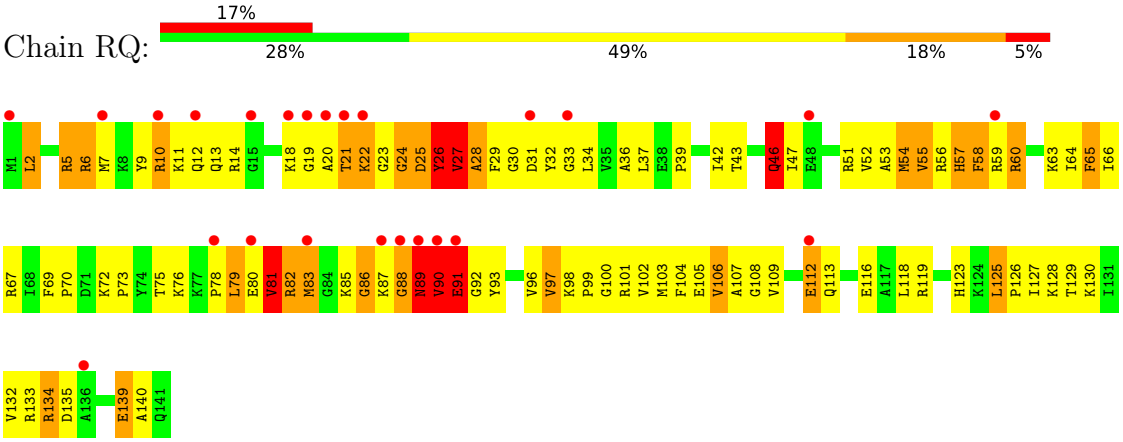
• Molecule 35: 50S ribosomal protein L15

Chain YP:



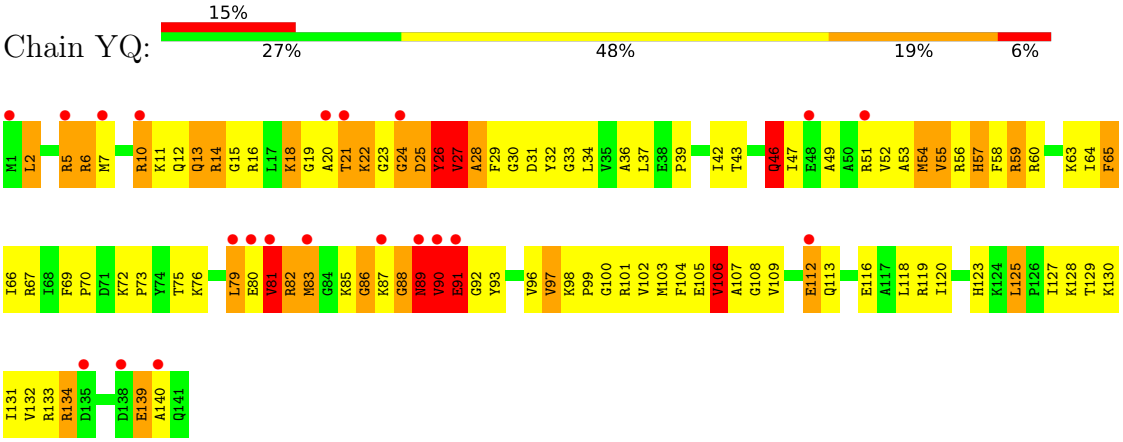
• Molecule 36: 50S ribosomal protein L16

Chain RQ:



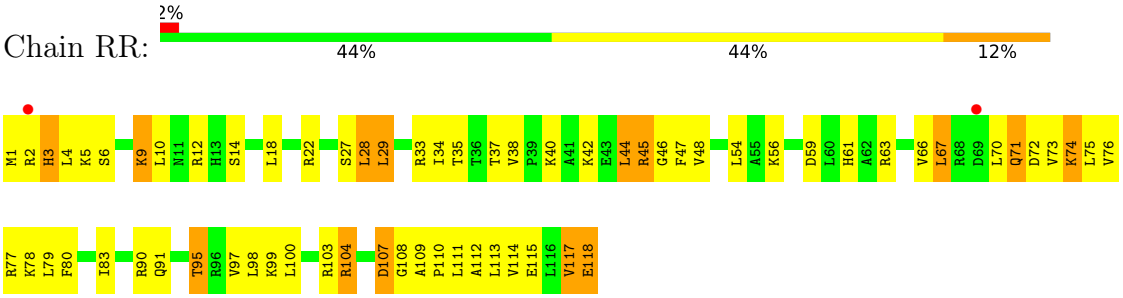
• Molecule 36: 50S ribosomal protein L16

Chain YQ:



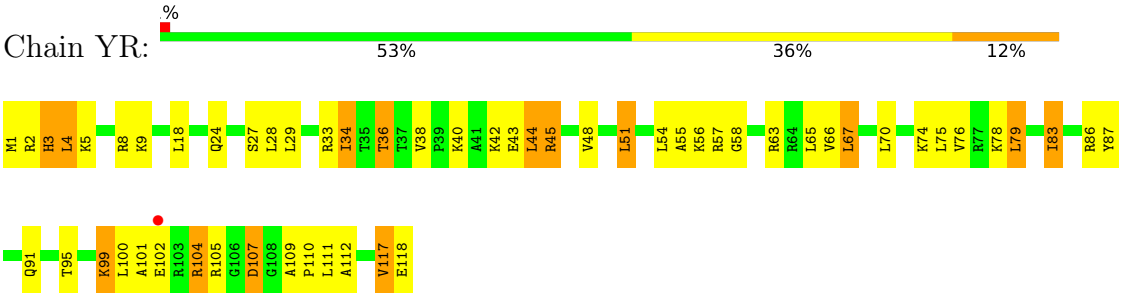
• Molecule 37: 50S ribosomal protein L17

Chain RR:



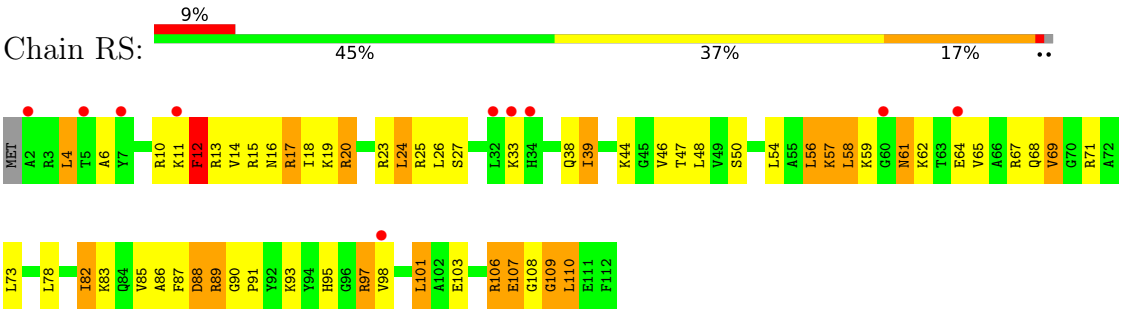
• Molecule 37: 50S ribosomal protein L17

Chain YR:



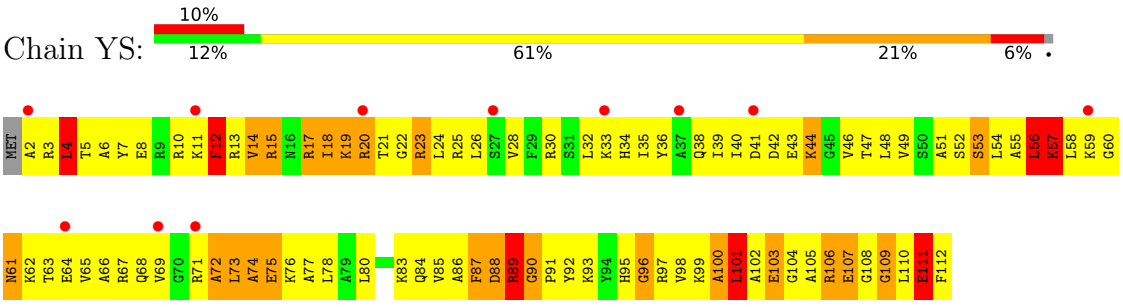
• Molecule 38: 50S ribosomal protein L18

Chain RS:



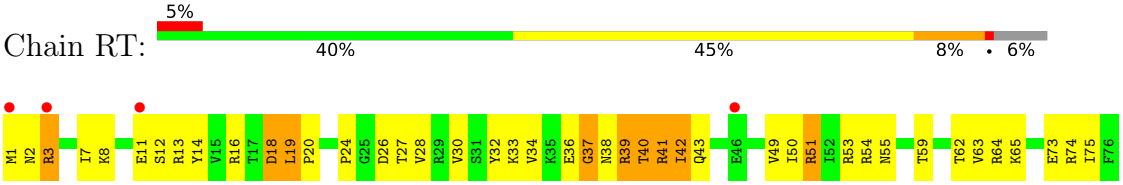
• Molecule 38: 50S ribosomal protein L18

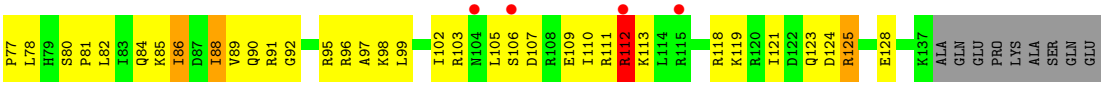
Chain YS:



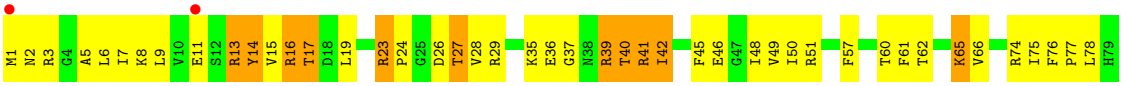
• Molecule 39: 50S ribosomal protein L19

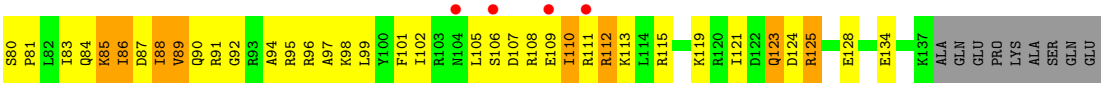
Chain RT:



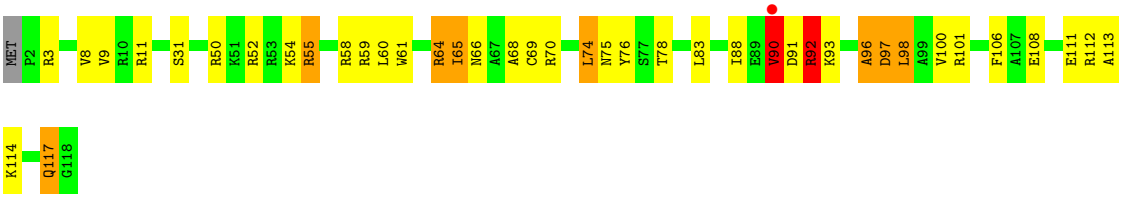


• Molecule 39: 50S ribosomal protein L19

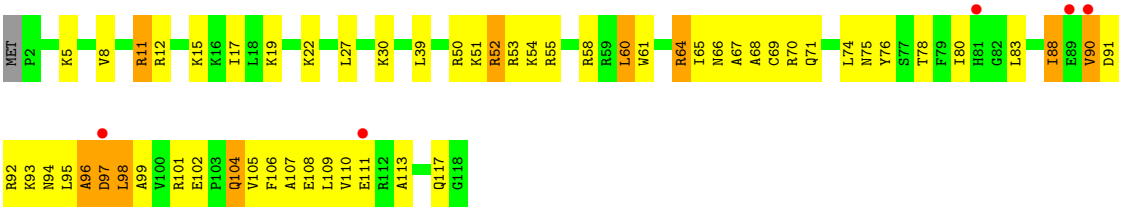




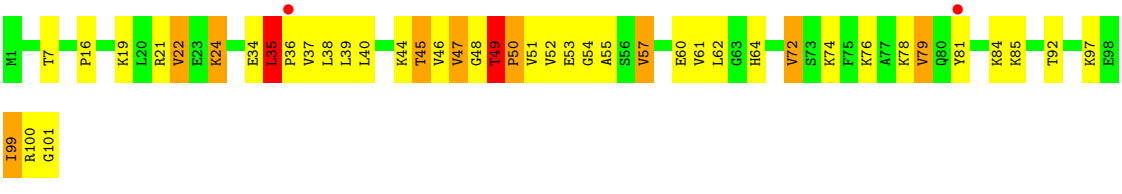
• Molecule 40: 50S ribosomal protein L20



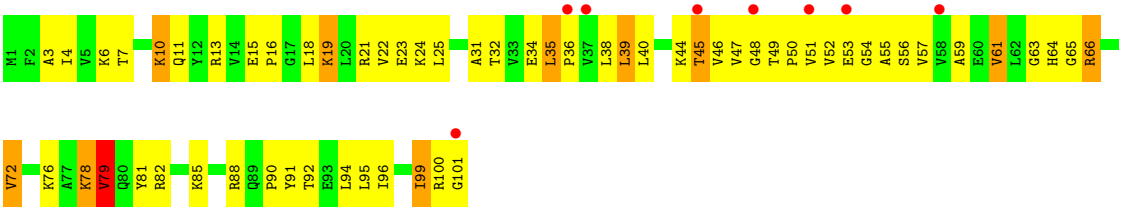
• Molecule 40: 50S ribosomal protein L20



• Molecule 41: 50S ribosomal protein L21

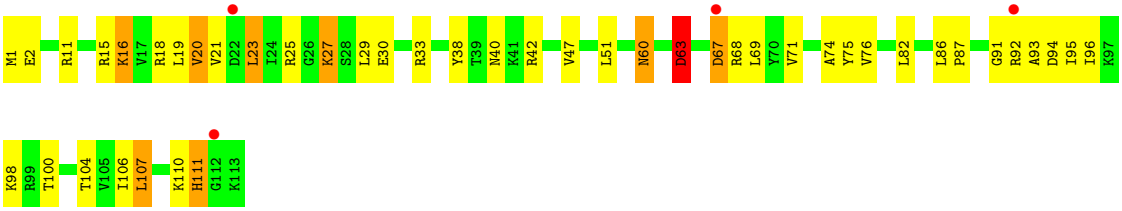


• Molecule 41: 50S ribosomal protein L21



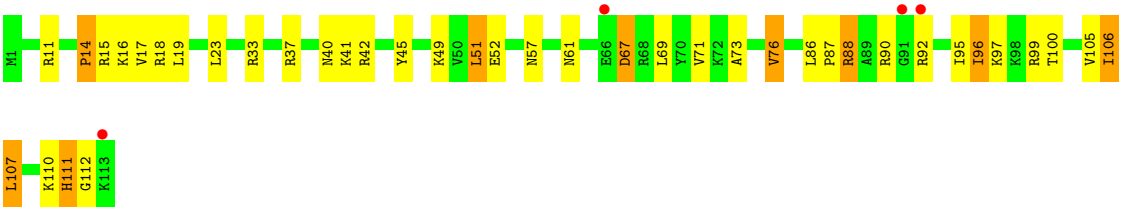
• Molecule 42: 50S ribosomal protein L22

Chain RW:



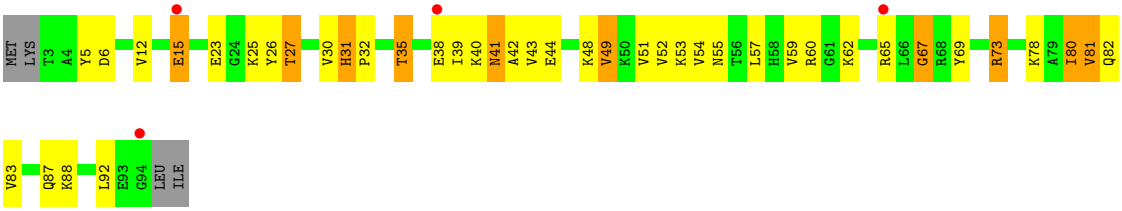
• Molecule 42: 50S ribosomal protein L22

Chain YW:



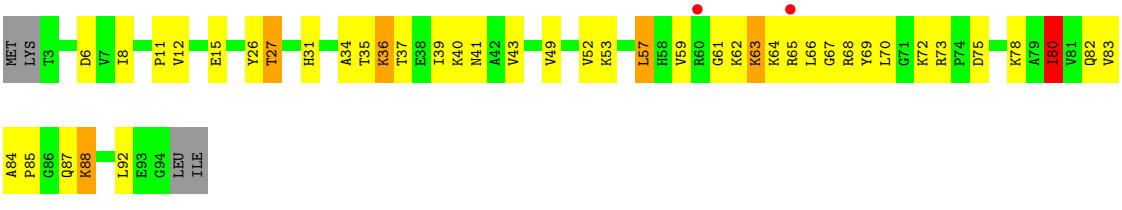
• Molecule 43: 50S ribosomal protein L23

Chain RX:



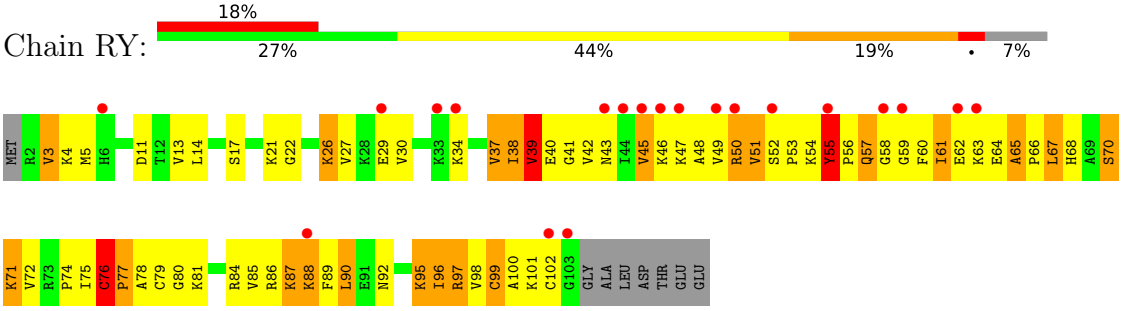
• Molecule 43: 50S ribosomal protein L23

Chain YX:



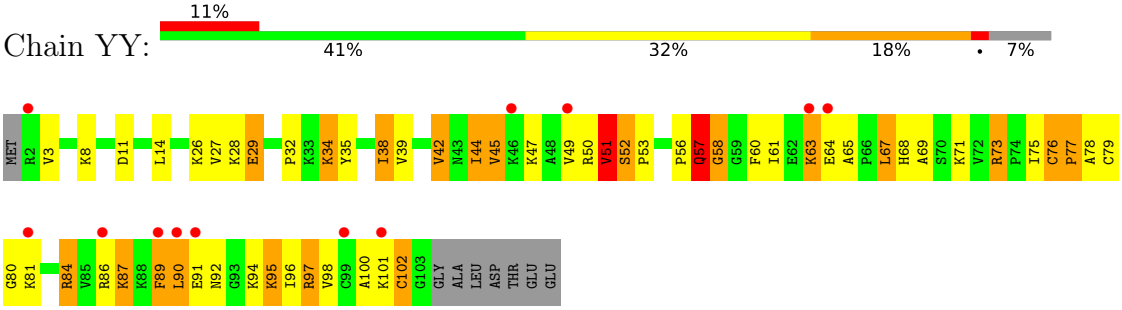
• Molecule 44: 50S ribosomal protein L24

Chain RY:



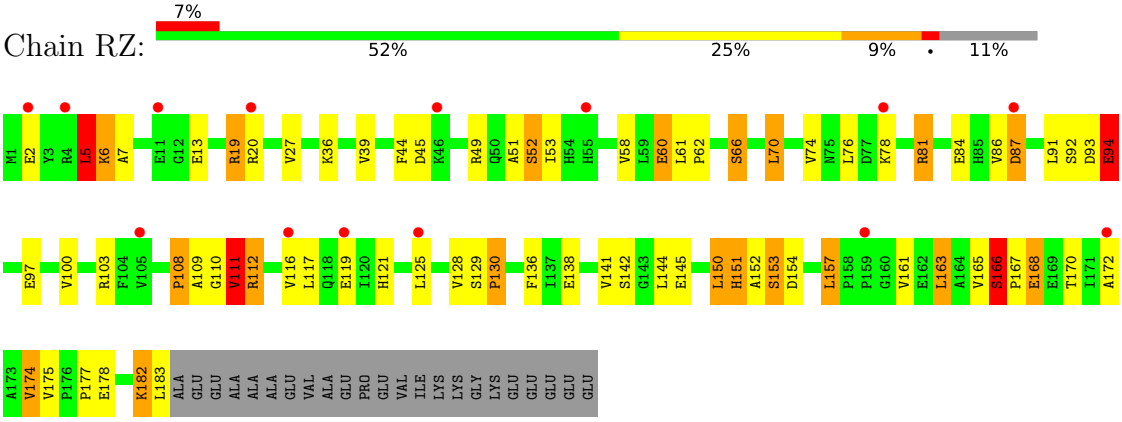
• Molecule 44: 50S ribosomal protein L24

Chain YY:



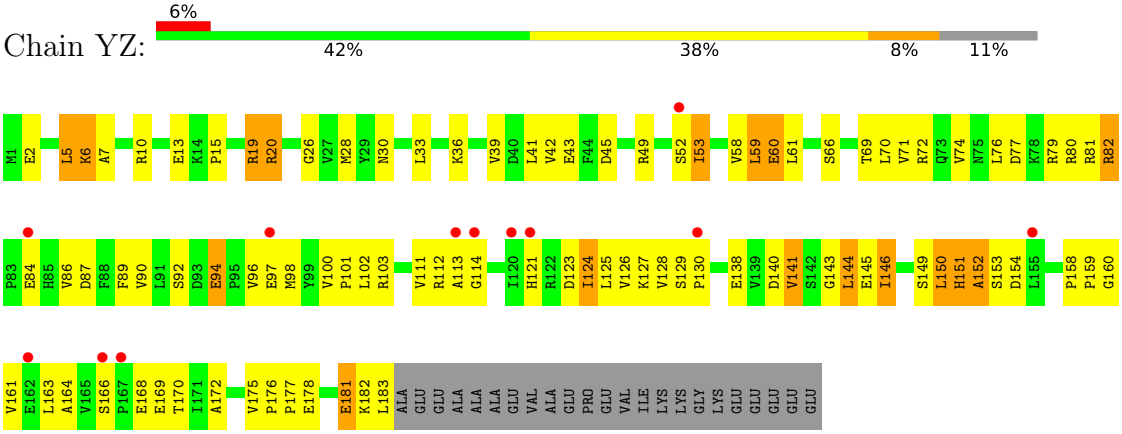
• Molecule 45: 50S ribosomal protein L25

Chain RZ:



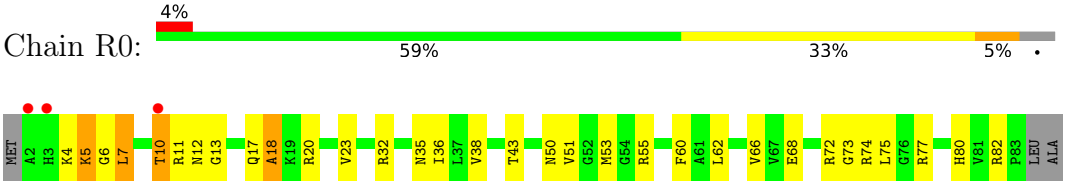
• Molecule 45: 50S ribosomal protein L25

Chain YZ:



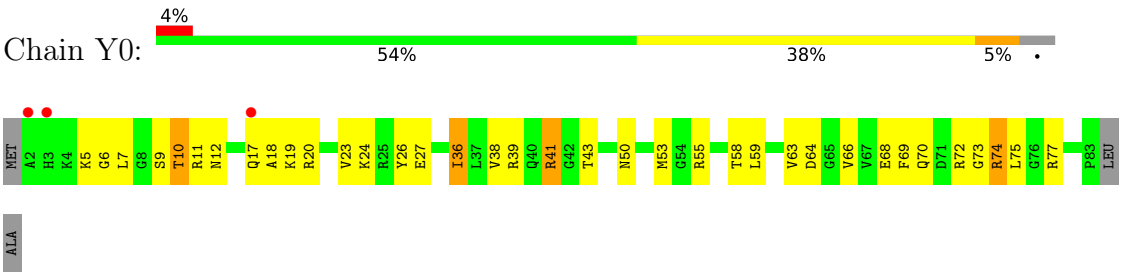
• Molecule 46: 50S ribosomal protein L27

Chain R0:



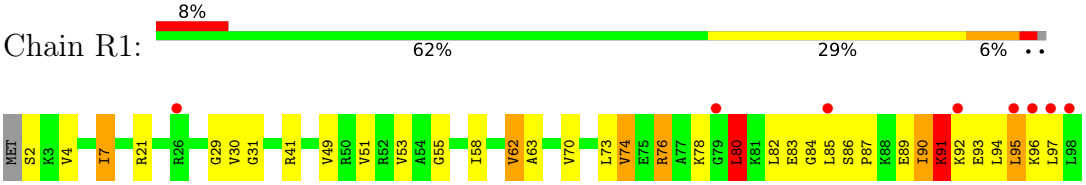
• Molecule 46: 50S ribosomal protein L27

Chain Y0:



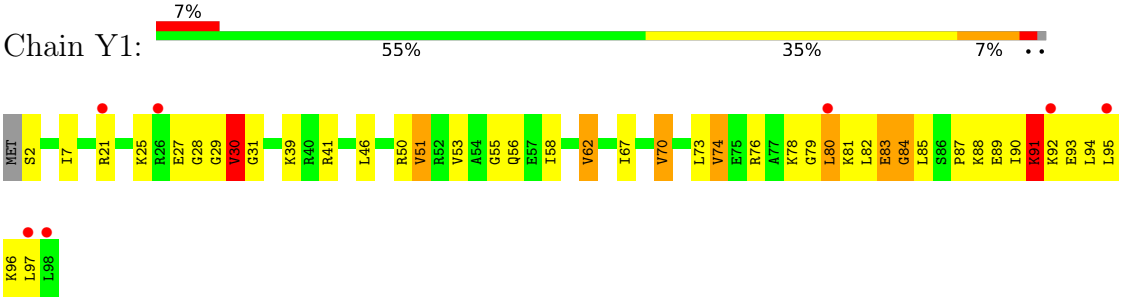
• Molecule 47: 50S ribosomal protein L28

Chain R1:



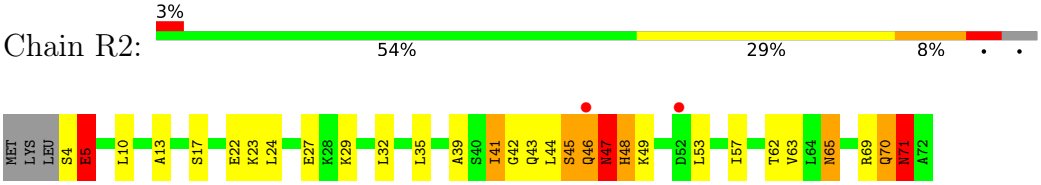
• Molecule 47: 50S ribosomal protein L28

Chain Y1:



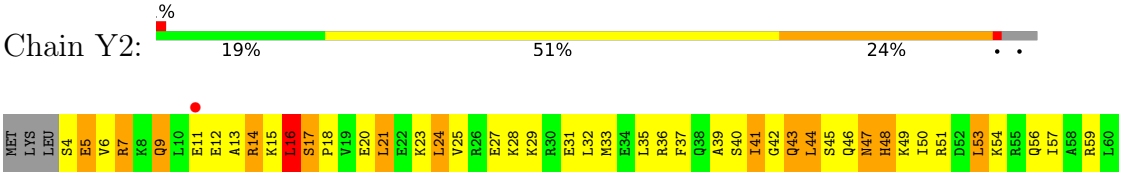
• Molecule 48: 50S ribosomal protein L29

Chain R2:



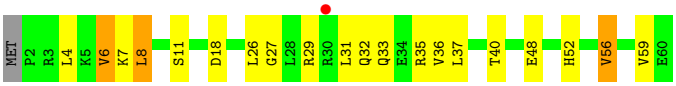
• Molecule 48: 50S ribosomal protein L29

Chain Y2:



• Molecule 49: 50S ribosomal protein L30

Chain R3:

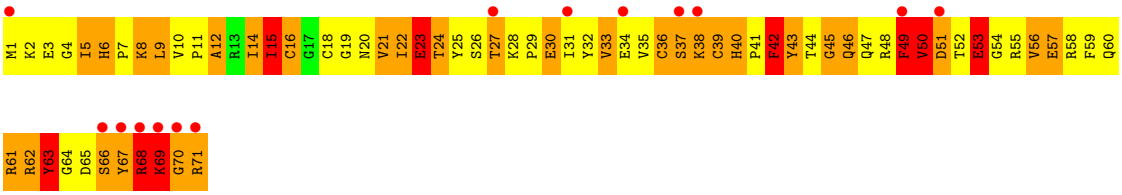


• Molecule 49: 50S ribosomal protein L30

Chain Y3:

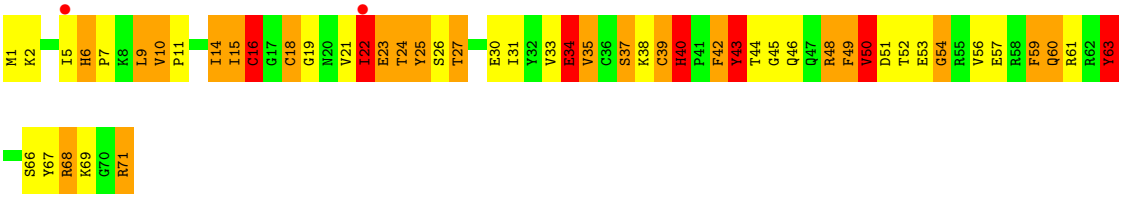
• Molecule 50: 50S ribosomal protein L31

Chain R4:



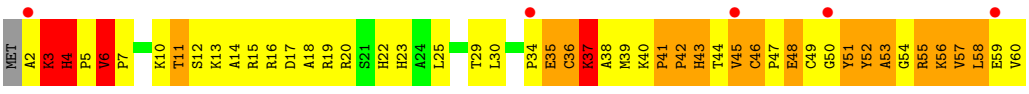
• Molecule 50: 50S ribosomal protein L31

Chain Y4:



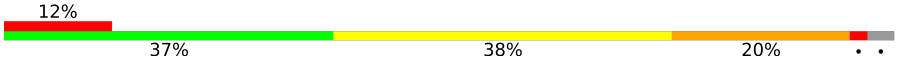
• Molecule 51: 50S ribosomal protein L32

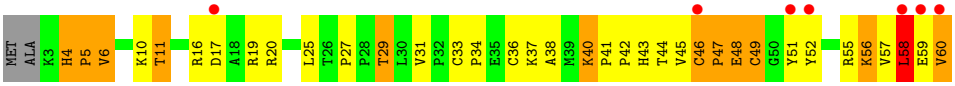
Chain R5:



• Molecule 51: 50S ribosomal protein L32

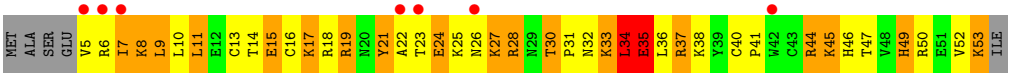
Chain Y5:





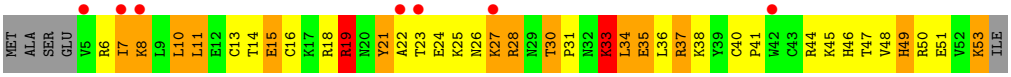
• Molecule 52: 50S ribosomal protein L33

Chain R6:



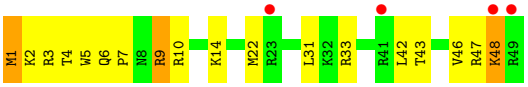
• Molecule 52: 50S ribosomal protein L33

Chain Y6:



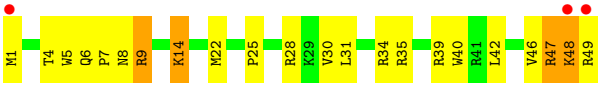
• Molecule 53: 50S ribosomal protein L34

Chain R7:



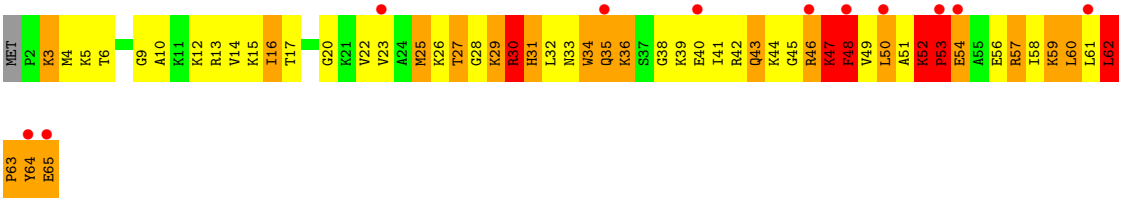
• Molecule 53: 50S ribosomal protein L34

Chain Y7:



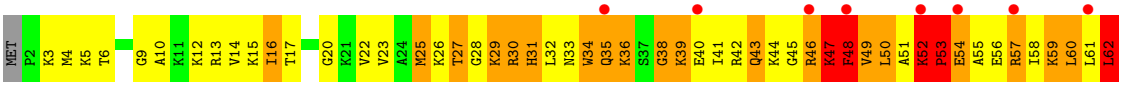
• Molecule 54: 50S ribosomal protein L35

Chain R8:



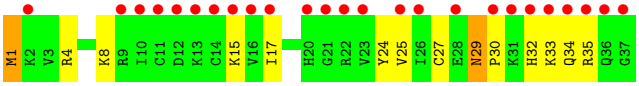
• Molecule 54: 50S ribosomal protein L35

Chain Y8:



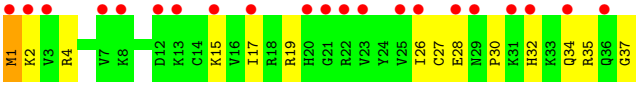
• Molecule 55: 50S ribosomal protein L36

Chain R9:



• Molecule 55: 50S ribosomal protein L36

Chain Y9:



• Molecule 56: CC-Puro

Chain Z5:

• Molecule 56: CC-Puro

Chain Z6:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	208.98Å 448.71Å 621.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.59 – 3.60 68.59 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.2 (68.59-3.60) 98.2 (68.59-3.60)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.46 (at 3.41Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, R_{free}	0.207 , 0.252 0.210 , 0.252	Depositor DCC
R_{free} test set	35866 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å^2)	80.8	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 79.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.42$, $< L^2 > = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	292311	wwPDB-VP
Average B, all atoms (Å^2)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.70% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: PAR, ZN, PPU, 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	QA	0.23	0/36097	0.44	2/56339 (0.0%)
1	XA	0.25	0/36101	0.45	3/56346 (0.0%)
2	QB	0.40	0/1959	0.87	4/2642 (0.2%)
2	XB	0.40	0/1959	0.89	4/2642 (0.2%)
3	QC	0.44	0/1629	0.87	0/2195
3	XC	0.48	0/1629	0.85	0/2195
4	QD	0.34	0/1733	0.81	4/2318 (0.2%)
4	XD	0.51	0/1733	0.96	5/2318 (0.2%)
5	QE	0.50	1/1171 (0.1%)	0.98	6/1576 (0.4%)
5	XE	0.54	0/1171	0.97	3/1576 (0.2%)
6	QF	0.45	0/856	0.80	0/1154
6	XF	0.51	0/856	0.87	0/1154
7	QG	0.43	0/1276	0.84	0/1709
7	XG	0.46	0/1276	0.88	2/1709 (0.1%)
8	QH	0.43	0/1136	0.93	2/1527 (0.1%)
8	XH	0.47	0/1136	0.92	4/1527 (0.3%)
9	QI	0.41	0/1029	0.90	1/1379 (0.1%)
9	XI	0.42	0/1029	0.89	1/1379 (0.1%)
10	QJ	0.55	1/814 (0.1%)	1.00	7/1095 (0.6%)
10	XJ	0.52	2/814 (0.2%)	1.03	6/1095 (0.5%)
11	QK	0.50	0/900	0.94	3/1213 (0.2%)
11	XK	0.51	0/900	0.93	0/1213
12	QL	0.59	1/991 (0.1%)	1.23	9/1327 (0.7%)
12	XL	0.59	0/991	1.26	12/1327 (0.9%)
13	QM	0.42	0/974	0.97	0/1303
13	XM	0.47	0/974	0.97	1/1303 (0.1%)
14	QN	0.49	0/501	0.96	1/664 (0.2%)
14	XN	0.56	0/501	1.04	3/664 (0.5%)
15	QO	0.45	0/745	0.80	0/992
15	XO	0.50	0/745	0.89	0/992
16	QP	0.48	0/721	0.93	2/970 (0.2%)
16	XP	0.48	0/721	0.88	1/970 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
17	QQ	0.42	0/847	0.85	0/1131
17	XQ	0.43	0/847	0.83	0/1131
18	QR	0.44	0/579	0.89	0/768
18	XR	0.47	0/579	0.85	1/768 (0.1%)
19	QS	0.45	0/689	0.99	3/926 (0.3%)
19	XS	0.47	0/689	1.01	6/926 (0.6%)
20	QT	0.47	0/765	0.96	2/1007 (0.2%)
20	XT	0.52	1/765 (0.1%)	0.97	3/1007 (0.3%)
21	QU	0.37	0/221	0.79	0/288
21	XU	0.37	0/221	0.81	0/288
22	QV	0.17	0/1836	0.34	0/2859
22	XV	0.18	0/1836	0.35	0/2859
23	QX	0.22	0/246	0.34	0/381
23	XX	0.43	0/246	0.60	0/381
24	QY	0.22	0/358	0.34	0/556
24	XY	0.33	0/384	0.46	0/597
25	RA	0.21	0/69521	0.40	3/108529 (0.0%)
25	YA	0.23	0/69521	0.42	3/108529 (0.0%)
26	RB	0.21	0/2878	0.39	0/4490
26	YB	0.23	0/2878	0.41	0/4490
27	RD	0.64	0/2165	1.07	10/2919 (0.3%)
27	YD	0.72	1/2165 (0.0%)	1.29	31/2919 (1.1%)
28	RE	0.62	0/1601	1.35	29/2160 (1.3%)
28	YE	0.62	0/1601	1.35	29/2160 (1.3%)
29	RF	0.38	0/1620	0.83	2/2194 (0.1%)
29	YF	0.62	0/1620	1.19	19/2194 (0.9%)
30	RG	0.41	0/1499	0.91	0/2016
30	YG	0.50	0/1499	0.91	2/2016 (0.1%)
31	RH	0.58	0/1332	1.22	14/1802 (0.8%)
31	YH	0.57	0/1332	1.22	15/1802 (0.8%)
32	RI	0.40	0/1151	0.89	2/1558 (0.1%)
32	YI	0.51	1/1151 (0.1%)	0.89	3/1558 (0.2%)
33	RN	0.52	0/1131	1.01	2/1525 (0.1%)
33	YN	0.55	0/1131	0.94	1/1525 (0.1%)
34	RO	0.55	0/943	0.98	3/1269 (0.2%)
34	YO	0.60	0/943	0.97	0/1269
35	RP	0.40	0/1162	0.90	2/1544 (0.1%)
35	YP	0.44	0/1162	0.96	2/1544 (0.1%)
36	RQ	0.71	0/1143	1.36	19/1527 (1.2%)
36	YQ	0.71	0/1143	1.34	19/1527 (1.2%)
37	RR	0.54	0/982	0.99	1/1312 (0.1%)
37	YR	0.56	0/982	1.04	3/1312 (0.2%)
38	RS	0.43	0/892	0.88	0/1187

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
38	YS	0.53	0/892	1.14	7/1187 (0.6%)
39	RT	0.53	0/1155	0.98	3/1542 (0.2%)
39	YT	0.56	0/1155	1.04	7/1542 (0.5%)
40	RU	0.53	0/982	1.04	6/1306 (0.5%)
40	YU	0.65	0/982	1.07	2/1306 (0.2%)
41	RV	0.47	0/790	0.97	4/1057 (0.4%)
41	YV	0.58	0/790	1.00	2/1057 (0.2%)
42	RW	0.64	0/911	0.98	0/1220
42	YW	0.60	0/911	1.00	0/1220
43	RX	0.59	0/739	0.96	4/993 (0.4%)
43	YX	0.60	0/739	1.00	1/993 (0.1%)
44	RY	0.54	0/798	1.07	6/1064 (0.6%)
44	YY	0.59	0/798	1.09	5/1064 (0.5%)
45	RZ	0.39	0/1493	0.88	8/2026 (0.4%)
45	YZ	0.39	0/1493	0.93	4/2026 (0.2%)
46	R0	0.56	0/657	1.04	5/874 (0.6%)
46	Y0	0.60	0/657	0.98	1/874 (0.1%)
47	R1	0.55	0/770	0.97	0/1022
47	Y1	0.59	0/770	0.96	1/1022 (0.1%)
48	R2	0.47	0/583	1.00	2/771 (0.3%)
48	Y2	0.60	0/583	1.36	9/771 (1.2%)
49	R3	0.46	0/474	0.91	0/635
49	Y3	0.53	0/474	0.86	0/635
50	R4	0.47	0/594	1.24	10/795 (1.3%)
50	Y4	0.48	0/594	1.16	6/795 (0.8%)
51	R5	0.65	0/473	1.22	6/639 (0.9%)
51	Y5	0.68	0/468	1.15	4/632 (0.6%)
52	R6	0.53	0/431	1.01	2/575 (0.3%)
52	Y6	0.56	0/431	1.00	2/575 (0.3%)
53	R7	0.62	0/438	1.05	2/575 (0.3%)
53	Y7	0.69	0/438	1.02	3/575 (0.5%)
54	R8	0.81	0/525	1.42	10/691 (1.4%)
54	Y8	0.79	0/525	1.41	9/691 (1.3%)
55	R9	0.35	0/310	0.81	0/407
55	Y9	0.45	0/310	0.80	0/407
56	Z5	0.45	0/40	0.63	0/60
56	Z6	0.45	0/40	0.62	0/60
All	All	0.34	8/316537 (0.0%)	0.64	446/473247 (0.1%)

All (8) bond length outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
32	YI	145	VAL	C-N	10.08	1.47	1.33
10	QJ	59	SER	C-N	-8.55	1.26	1.33
20	XT	50	GLU	C-N	8.19	1.46	1.33
12	QL	48	PRO	N-CD	5.45	1.55	1.47
5	QE	70	PRO	N-CD	5.31	1.55	1.47
10	XJ	53	PRO	N-CD	5.22	1.55	1.47
10	XJ	52	GLY	C-N	5.10	1.41	1.34
27	YD	241	PRO	N-CD	5.02	1.54	1.47

All (446) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
50	R4	39	CYS	N-CA-C	-13.60	95.95	112.54
28	RE	58	ARG	N-CA-C	-11.79	95.27	113.89
28	YE	58	ARG	N-CA-C	-11.78	95.28	113.89
48	Y2	5	GLU	N-CA-C	-11.56	99.66	113.88
12	XL	119	LYS	N-CA-C	-10.39	95.14	110.52
36	RQ	89	ASN	N-CA-C	-10.38	100.42	112.87
12	QL	119	LYS	N-CA-C	-10.34	95.22	110.52
36	YQ	89	ASN	N-CA-C	-10.28	100.53	112.87
28	YE	21	VAL	C-N-CD	-10.10	98.38	120.60
28	RE	21	VAL	C-N-CD	-10.08	98.42	120.60
36	YQ	81	VAL	CB-CA-C	-10.02	94.86	111.29
36	YQ	81	VAL	N-CA-C	9.94	130.02	109.34
36	RQ	81	VAL	CB-CA-C	-9.91	95.03	111.29
36	RQ	81	VAL	N-CA-C	9.91	129.95	109.34
33	RN	114	ARG	N-CA-C	-9.75	97.95	110.53
50	Y4	39	CYS	N-CA-C	-9.44	101.31	112.92
12	QL	26	ALA	N-CA-C	-9.26	102.57	113.21
29	YF	197	ASP	N-CA-C	-9.22	98.76	110.19
12	XL	26	ALA	N-CA-C	-9.20	102.64	113.21
50	R4	45	GLY	N-CA-C	-9.13	102.72	115.32
12	XL	47	LYS	C-N-CD	9.04	140.50	120.60
40	YU	96	ALA	N-CA-C	-9.02	98.92	112.54
19	QS	6	LYS	N-CA-C	-9.01	102.68	114.31
5	XE	69	VAL	O-C-N	8.91	131.26	121.10
48	Y2	17	SER	C-N-CD	-8.83	101.18	120.60
27	YD	91	ARG	N-CA-C	-8.73	98.76	110.55
54	Y8	36	LYS	N-CA-C	-8.71	96.64	110.14
36	RQ	125	LEU	CA-C-N	8.69	130.70	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
36	RQ	125	LEU	C-N-CA	8.69	130.70	119.84
36	YQ	125	LEU	CA-C-N	8.69	130.70	119.84
36	YQ	125	LEU	C-N-CA	8.69	130.70	119.84
54	R8	36	LYS	N-CA-C	-8.63	96.76	110.14
35	YP	59	LEU	N-CA-C	-8.50	102.62	113.16
14	QN	26	ARG	N-CA-C	8.50	120.32	111.14
28	RE	23	VAL	N-CA-C	8.46	120.97	108.96
28	YE	23	VAL	N-CA-C	8.43	120.93	108.96
27	YD	193	VAL	N-CA-C	8.27	120.25	109.58
29	YF	75	HIS	N-CA-C	8.23	122.28	109.52
7	XG	114	ARG	N-CA-C	8.10	120.19	111.36
31	YH	19	VAL	N-CA-C	8.08	117.92	106.53
12	QL	47	LYS	C-N-CD	8.06	138.34	120.60
29	YF	65	TRP	CA-C-N	8.05	129.91	119.84
29	YF	65	TRP	C-N-CA	8.05	129.91	119.84
19	QS	8	GLY	N-CA-C	8.04	119.57	111.21
31	RH	19	VAL	N-CA-C	8.03	117.85	106.53
50	R4	63	TYR	N-CA-C	-7.99	102.80	112.54
54	Y8	45	GLY	N-CA-C	-7.99	101.79	112.81
54	R8	45	GLY	N-CA-C	-7.95	101.84	112.81
50	Y4	63	TYR	N-CA-C	-7.87	103.82	113.50
36	YQ	5	ARG	N-CA-C	-7.76	96.93	109.96
51	R5	6	VAL	CA-C-N	7.74	128.96	119.98
51	R5	6	VAL	C-N-CA	7.74	128.96	119.98
31	YH	100	GLY	N-CA-C	-7.73	99.72	111.24
31	RH	100	GLY	N-CA-C	-7.71	99.76	111.24
36	RQ	5	ARG	N-CA-C	-7.70	97.02	109.96
11	QK	40	ILE	N-CA-C	-7.70	105.66	111.90
37	YR	9	LYS	N-CA-C	-7.64	98.17	110.32
10	QJ	59	SER	O-C-N	7.63	132.74	122.59
48	Y2	14	ARG	N-CA-C	-7.62	99.98	110.35
54	R8	38	GLY	N-CA-C	-7.62	103.16	113.37
27	YD	251	GLY	N-CA-C	7.61	126.75	115.64
54	Y8	38	GLY	N-CA-C	-7.61	103.18	113.37
53	Y7	6	GLN	CA-C-N	7.59	127.63	119.89
53	Y7	6	GLN	C-N-CA	7.59	127.63	119.89
28	RE	185	LYS	N-CA-C	7.59	119.94	109.54
31	RH	157	TYR	N-CA-C	-7.58	103.10	112.88
31	YH	157	TYR	N-CA-C	-7.56	103.12	112.88
48	Y2	21	LEU	N-CA-C	-7.54	103.00	111.07
12	XL	58	VAL	N-CA-C	7.51	118.89	107.77
4	XD	31	CYS	N-CA-C	-7.51	104.16	112.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
12	QL	58	VAL	N-CA-C	7.49	118.86	107.77
28	YE	185	LYS	N-CA-C	7.49	119.81	109.54
27	YD	43	ARG	N-CA-C	7.49	122.06	113.15
48	Y2	41	ILE	N-CA-C	-7.48	103.26	112.76
31	RH	117	PRO	N-CA-C	7.43	119.77	110.70
31	YH	117	PRO	N-CA-C	7.37	119.69	110.70
12	QL	44	THR	CA-C-N	7.36	129.04	119.84
12	QL	44	THR	C-N-CA	7.36	129.04	119.84
27	RD	244	ARG	CA-C-N	-7.33	112.83	120.38
27	RD	244	ARG	C-N-CA	-7.33	112.83	120.38
12	XL	44	THR	CA-C-N	7.31	128.98	119.84
12	XL	44	THR	C-N-CA	7.31	128.98	119.84
54	R8	54	GLU	N-CA-C	-7.29	102.52	111.33
4	QD	22	LYS	N-CA-C	7.25	119.26	111.36
31	RH	139	GLN	N-CA-C	-7.24	103.39	111.28
54	Y8	54	GLU	N-CA-C	-7.23	102.58	111.33
48	Y2	16	LEU	N-CA-C	-7.21	95.45	110.80
27	RD	272	ALA	N-CA-C	7.19	118.15	108.23
31	YH	139	GLN	N-CA-C	-7.14	103.49	111.28
27	YD	53	PHE	N-CA-C	7.13	121.56	112.86
53	R7	6	GLN	CA-C-N	7.11	127.15	119.90
53	R7	6	GLN	C-N-CA	7.11	127.15	119.90
28	RE	168	MET	N-CA-C	7.10	121.09	109.94
28	YE	168	MET	N-CA-C	7.09	121.07	109.94
34	RO	100	GLY	CA-C-N	7.06	127.02	120.03
34	RO	100	GLY	C-N-CA	7.06	127.02	120.03
29	YF	203	GLN	N-CA-C	-7.06	103.28	110.97
33	RN	85	ILE	N-CA-C	7.00	115.71	107.73
16	XP	20	VAL	N-CA-C	6.98	118.40	108.42
28	YE	188	VAL	CA-C-N	6.96	128.54	119.84
28	YE	188	VAL	C-N-CA	6.96	128.54	119.84
45	YZ	158	PRO	CA-C-N	6.94	128.52	119.84
45	YZ	158	PRO	C-N-CA	6.94	128.52	119.84
11	QK	102	GLY	N-CA-C	-6.91	96.81	113.18
28	RE	188	VAL	CA-C-N	6.90	128.47	119.84
28	RE	188	VAL	C-N-CA	6.90	128.47	119.84
36	YQ	116	GLU	N-CA-C	-6.89	103.21	111.69
32	YI	133	HIS	CA-C-N	-6.86	112.59	119.93
32	YI	133	HIS	C-N-CA	-6.86	112.59	119.93
50	Y4	59	PHE	N-CA-C	-6.85	103.56	112.68
36	YQ	10	ARG	N-CA-C	-6.84	99.16	110.17
36	RQ	10	ARG	N-CA-C	-6.84	99.16	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
29	YF	24	LEU	CA-C-N	6.83	128.38	119.84
29	YF	24	LEU	C-N-CA	6.83	128.38	119.84
40	RU	76	TYR	N-CA-C	-6.82	103.93	111.36
35	RP	59	LEU	N-CA-C	-6.80	104.56	112.92
36	RQ	116	GLU	N-CA-C	-6.77	103.36	111.69
5	QE	23	GLY	N-CA-C	6.76	120.58	110.97
40	RU	96	ALA	N-CA-C	-6.75	103.35	112.26
36	YQ	46	GLN	N-CA-C	-6.71	105.10	113.55
20	QT	12	ALA	N-CA-C	-6.70	102.42	112.54
36	RQ	46	GLN	N-CA-C	-6.70	105.11	113.55
27	YD	223	GLY	N-CA-C	-6.68	104.41	112.49
27	YD	198	ASN	N-CA-C	-6.67	102.28	111.28
46	R0	17	GLN	N-CA-C	6.64	119.72	108.90
2	QB	166	ASP	CA-C-N	6.63	126.25	119.56
2	QB	166	ASP	C-N-CA	6.63	126.25	119.56
40	RU	97	ASP	N-CA-C	-6.62	103.72	112.72
36	YQ	65	PHE	N-CA-C	6.62	121.89	113.55
20	XT	12	ALA	N-CA-C	-6.60	104.85	113.17
38	YS	15	ARG	N-CA-C	-6.60	103.57	111.69
28	YE	95	ILE	N-CA-C	-6.57	104.46	113.00
36	RQ	65	PHE	N-CA-C	6.56	121.81	113.55
27	YD	14	ARG	N-CA-C	-6.56	104.33	112.72
50	Y4	40	HIS	CA-C-N	-6.55	113.03	119.78
50	Y4	40	HIS	C-N-CA	-6.55	113.03	119.78
50	Y4	45	GLY	N-CA-C	-6.55	106.19	115.30
28	RE	95	ILE	N-CA-C	-6.55	104.48	113.00
25	YA	1360	A	C2'-C3'-O3'	-6.54	103.88	113.70
36	RQ	82	ARG	N-CA-C	6.54	123.83	114.16
51	Y5	46	CYS	CA-C-N	6.53	128.00	119.84
51	Y5	46	CYS	C-N-CA	6.53	128.00	119.84
36	YQ	82	ARG	N-CA-C	6.51	123.79	114.16
31	YH	164	TYR	N-CA-C	-6.49	100.68	110.28
19	QS	67	VAL	N-CA-C	6.48	117.27	110.72
31	RH	164	TYR	N-CA-C	-6.46	100.72	110.28
28	YE	125	GLY	CA-C-N	6.46	127.92	119.84
28	YE	125	GLY	C-N-CA	6.46	127.92	119.84
12	XL	47	LYS	CB-CA-C	6.44	118.55	108.63
54	Y8	48	PHE	N-CA-C	6.43	118.89	108.34
28	RE	125	GLY	CA-C-N	6.41	127.86	119.84
28	RE	125	GLY	C-N-CA	6.41	127.86	119.84
39	YT	111	ARG	N-CA-C	-6.41	102.53	110.41
44	RY	65	ALA	CA-C-N	6.41	126.42	119.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
44	RY	65	ALA	C-N-CA	6.41	126.42	119.89
54	R8	48	PHE	N-CA-C	6.40	118.84	108.34
54	R8	50	LEU	N-CA-C	6.40	118.92	108.55
27	YD	83	GLU	N-CA-C	6.39	118.89	109.69
45	RZ	157	LEU	CA-C-N	6.37	126.94	120.38
45	RZ	157	LEU	C-N-CA	6.37	126.94	120.38
40	YU	97	ASP	N-CA-C	-6.37	105.49	113.20
32	RI	135	GLU	N-CA-C	6.35	116.83	108.07
16	QP	71	ARG	N-CA-C	-6.34	104.45	111.36
4	QD	171	GLY	CA-C-N	6.33	126.02	119.56
4	QD	171	GLY	C-N-CA	6.33	126.02	119.56
54	Y8	50	LEU	N-CA-C	6.32	118.78	108.55
29	RF	65	TRP	CA-C-N	6.31	127.73	119.84
29	RF	65	TRP	C-N-CA	6.31	127.73	119.84
52	R6	24	GLU	N-CA-C	6.31	116.33	108.19
54	R8	23	VAL	N-CA-C	6.30	117.09	107.77
35	RP	25	SER	N-CA-C	-6.28	104.88	113.30
36	RQ	106	VAL	N-CA-C	6.28	116.19	108.53
40	RU	9	VAL	N-CA-C	6.27	116.44	110.42
54	Y8	23	VAL	N-CA-C	6.27	117.04	107.77
36	YQ	106	VAL	N-CA-C	6.25	116.16	108.53
41	YV	35	LEU	CA-C-N	6.25	127.65	119.84
41	YV	35	LEU	C-N-CA	6.25	127.65	119.84
12	XL	89	ARG	N-CA-C	6.23	118.57	110.53
27	YD	272	ALA	N-CA-C	6.22	124.06	110.80
40	RU	92	ARG	N-CA-C	-6.21	104.29	112.24
50	R4	68	ARG	N-CA-C	6.19	123.99	110.80
12	XL	47	LYS	N-CA-C	-6.19	101.55	110.40
27	YD	226	MET	N-CA-C	6.17	117.83	110.19
27	YD	22	SER	N-CA-C	-6.16	103.89	111.40
40	RU	74	LEU	N-CA-C	6.15	116.36	108.24
43	YX	80	ILE	CB-CA-C	6.15	118.98	110.99
31	RH	127	GLU	N-CA-C	-6.15	96.22	109.81
31	YH	127	GLU	N-CA-C	-6.15	96.22	109.81
39	YT	57	PHE	N-CA-C	-6.15	102.85	110.41
36	YQ	96	VAL	N-CA-C	6.14	116.09	107.37
27	YD	69	ARG	N-CA-C	6.13	120.76	113.28
39	YT	14	TYR	N-CA-C	-6.12	105.86	113.15
46	R0	60	PHE	N-CA-C	6.11	118.46	109.24
36	RQ	97	VAL	N-CA-C	6.10	117.09	108.36
43	RX	31	HIS	CA-C-N	6.10	125.99	119.28
43	RX	31	HIS	C-N-CA	6.10	125.99	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	XB	25	ASN	CA-C-N	6.09	127.46	119.84
2	XB	25	ASN	C-N-CA	6.09	127.46	119.84
12	QL	89	ARG	N-CA-C	6.09	118.38	110.53
27	YD	107	ALA	N-CA-C	6.08	118.42	109.42
27	YD	230	ASP	N-CA-C	6.07	119.93	111.56
27	RD	49	ILE	N-CA-CB	6.07	117.14	110.53
36	RQ	96	VAL	N-CA-C	6.07	115.98	107.37
52	R6	28	ARG	N-CA-C	6.06	118.39	109.24
48	Y2	45	SER	N-CA-C	6.06	118.28	109.07
28	YE	77	ILE	N-CA-C	-6.05	103.57	112.04
30	YG	31	VAL	CA-C-N	6.02	127.36	119.84
30	YG	31	VAL	C-N-CA	6.02	127.36	119.84
28	RE	77	ILE	N-CA-C	-6.01	103.63	112.04
37	RR	9	LYS	N-CA-C	-6.00	100.78	109.59
28	YE	18	ASP	N-CA-C	-6.00	98.95	109.02
27	YD	88	ARG	N-CA-C	5.99	117.35	108.60
36	YQ	97	VAL	N-CA-C	5.99	116.93	108.36
41	RV	35	LEU	CA-C-N	5.97	127.31	119.84
41	RV	35	LEU	C-N-CA	5.97	127.31	119.84
54	R8	60	LEU	N-CA-C	-5.95	104.16	112.25
28	RE	18	ASP	N-CA-C	-5.94	99.03	109.02
39	RT	19	LEU	CA-C-N	5.94	125.85	119.85
39	RT	19	LEU	C-N-CA	5.94	125.85	119.85
29	YF	183	VAL	N-CA-C	5.94	117.39	110.62
50	R4	56	VAL	N-CA-C	-5.94	108.07	113.71
32	YI	53	ALA	O-C-N	5.94	128.44	122.03
54	Y8	60	LEU	N-CA-C	-5.93	104.18	112.25
31	YH	4	ILE	N-CA-C	5.93	117.31	108.23
54	R8	27	THR	N-CA-C	-5.93	102.14	110.50
38	YS	111	GLU	N-CA-C	-5.93	103.97	110.91
44	YY	11	ASP	N-CA-C	5.90	117.91	110.24
27	YD	196	VAL	N-CA-C	5.90	120.39	111.89
20	XT	50	GLU	O-C-N	-5.89	115.02	122.27
31	YH	125	VAL	C-N-CD	-5.89	107.64	120.60
31	RH	125	VAL	C-N-CD	-5.89	107.64	120.60
36	RQ	81	VAL	CA-C-N	-5.89	111.10	120.70
36	RQ	81	VAL	C-N-CA	-5.89	111.10	120.70
27	YD	106	ILE	N-CA-C	-5.89	101.94	109.30
31	RH	4	ILE	N-CA-C	5.88	117.23	108.23
5	QE	127	ASN	CA-C-N	5.88	126.02	119.32
5	QE	127	ASN	C-N-CA	5.88	126.02	119.32
39	YT	89	VAL	N-CA-C	-5.88	107.02	113.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
51	R5	4	HIS	CA-C-N	-5.87	113.89	119.76
51	R5	4	HIS	C-N-CA	-5.87	113.89	119.76
27	YD	200	ASP	N-CA-C	-5.87	105.61	112.89
14	XN	13	THR	CA-C-N	5.86	127.17	119.84
14	XN	13	THR	C-N-CA	5.86	127.17	119.84
19	XS	41	VAL	CA-C-N	5.86	141.06	127.00
19	XS	41	VAL	C-N-CA	5.86	141.06	127.00
51	Y5	58	LEU	N-CA-C	5.84	118.70	110.23
36	YQ	81	VAL	CA-C-N	-5.84	111.19	120.70
36	YQ	81	VAL	C-N-CA	-5.84	111.19	120.70
41	RV	49	THR	CA-C-N	-5.83	112.56	119.84
41	RV	49	THR	C-N-CA	-5.83	112.56	119.84
29	YF	151	SER	N-CA-C	-5.82	106.16	113.20
27	RD	13	ARG	N-CA-C	-5.81	106.23	113.38
28	RE	118	LYS	N-CA-C	-5.80	98.44	110.80
44	YY	89	PHE	N-CA-C	-5.80	104.37	113.02
31	RH	161	GLY	N-CA-C	5.80	118.30	111.63
38	YS	12	PHE	N-CA-C	-5.79	98.46	110.80
28	YE	118	LYS	N-CA-C	-5.79	98.47	110.80
46	R0	4	LYS	CA-C-N	-5.79	112.01	120.75
46	R0	4	LYS	C-N-CA	-5.79	112.01	120.75
29	YF	85	GLY	N-CA-C	-5.79	107.67	115.36
12	XL	48	PRO	CA-N-CD	-5.75	103.45	111.50
5	QE	69	VAL	O-C-N	5.75	126.91	121.11
11	QK	98	LEU	N-CA-C	5.74	117.34	111.14
50	R4	67	TYR	N-CA-C	5.74	118.56	109.50
31	YH	161	GLY	N-CA-C	5.74	118.23	111.63
27	RD	44	ASN	N-CA-C	5.71	116.83	108.14
45	RZ	178	GLU	N-CA-C	5.71	115.64	108.45
10	XJ	75	ILE	N-CA-C	5.70	115.89	110.53
28	YE	159	HIS	N-CA-C	-5.70	101.97	110.23
38	YS	6	ALA	N-CA-C	-5.69	105.17	111.71
43	RX	6	ASP	N-CA-C	-5.68	106.39	113.55
28	YE	131	ALA	CB-CA-C	-5.68	110.04	116.63
28	RE	49	LEU	N-CA-C	5.67	119.72	112.12
8	XH	26	VAL	CA-C-N	5.67	126.92	119.84
8	XH	26	VAL	C-N-CA	5.67	126.92	119.84
4	XD	39	PRO	CA-C-N	5.65	125.63	120.03
4	XD	39	PRO	C-N-CA	5.65	125.63	120.03
10	QJ	52	GLY	CA-C-N	5.65	125.01	118.85
10	QJ	52	GLY	C-N-CA	5.65	125.01	118.85
12	XL	46	LYS	N-CA-C	-5.64	100.97	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
28	YE	198	VAL	N-CA-C	5.63	116.23	107.28
28	RE	198	VAL	N-CA-C	5.63	116.23	107.28
38	YS	72	ALA	N-CA-C	-5.62	105.07	111.14
27	YD	11	PRO	N-CA-C	-5.62	102.18	111.11
25	YA	271(B)	G	P-O3'-C3'	5.61	126.44	119.70
8	XH	75	ARG	CA-C-N	5.61	126.85	119.84
8	XH	75	ARG	C-N-CA	5.61	126.85	119.84
28	RE	131	ALA	CB-CA-C	-5.61	110.12	116.63
28	YE	49	LEU	N-CA-C	5.61	119.63	112.12
28	YE	174	ASP	N-CA-C	5.60	115.63	108.24
46	Y0	55	ARG	N-CA-C	5.59	118.14	111.71
51	Y5	27	PRO	O-C-N	5.59	123.88	121.31
27	YD	268	ARG	N-CA-C	5.59	119.60	112.34
28	RE	93	VAL	N-CA-C	-5.59	97.72	109.34
29	YF	97	TYR	N-CA-C	5.59	119.57	112.87
27	YD	120	GLY	CA-C-N	5.58	125.42	119.28
27	YD	120	GLY	C-N-CA	5.58	125.42	119.28
28	RE	159	HIS	N-CA-C	-5.58	102.15	110.23
36	YQ	21	THR	N-CA-C	5.58	122.67	110.80
28	YE	93	VAL	N-CA-C	-5.57	97.75	109.34
10	QJ	55	LYS	CB-CA-C	-5.56	110.14	116.54
27	YD	174	ILE	N-CA-C	5.56	116.88	109.37
10	XJ	28	ARG	N-CA-C	-5.56	106.45	113.18
52	Y6	24	GLU	N-CA-C	5.56	115.45	108.45
14	XN	47	LEU	N-CA-C	5.55	117.33	111.28
46	R0	11	ARG	N-CA-C	-5.55	106.55	113.38
4	QD	31	CYS	N-CA-C	5.55	118.25	111.82
10	XJ	55	LYS	CB-CA-C	-5.55	110.16	116.54
36	RQ	21	THR	N-CA-C	5.54	122.61	110.80
36	YQ	26	TYR	N-CA-C	5.54	119.13	111.71
1	QA	189	U	P-O3'-C3'	5.53	128.49	120.20
50	R4	15	ILE	N-CA-C	5.53	118.07	108.90
9	XI	63	ILE	N-CA-C	5.53	116.01	107.78
28	YE	87	GLU	N-CA-C	-5.53	103.85	111.81
27	YD	146	GLU	N-CA-C	-5.52	101.86	110.42
31	RH	38	SER	CA-C-N	5.52	125.19	119.56
31	RH	38	SER	C-N-CA	5.52	125.19	119.56
9	QI	86	VAL	N-CA-C	-5.52	107.37	112.83
38	YS	53	SER	N-CA-C	-5.52	106.72	113.50
1	XA	812	C	C2'-C3'-O3'	5.51	117.77	109.50
12	QL	46	LYS	N-CA-C	-5.50	101.19	109.60
19	XS	58	VAL	N-CA-C	5.50	113.88	107.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
28	RE	87	GLU	N-CA-C	-5.48	103.92	111.81
29	YF	46	ARG	N-CA-C	5.47	119.67	112.89
12	XL	127	GLU	N-CA-C	-5.47	105.83	113.37
2	QB	25	ASN	CA-C-N	5.46	126.67	119.84
2	QB	25	ASN	C-N-CA	5.46	126.67	119.84
44	RY	76	CYS	CA-C-N	5.45	126.66	119.84
44	RY	76	CYS	C-N-CA	5.45	126.66	119.84
38	YS	101	LEU	N-CA-C	-5.45	104.99	111.69
33	YN	24	GLY	N-CA-C	-5.44	106.08	113.37
18	XR	25	THR	N-CA-C	-5.43	106.66	113.72
1	QA	189	U	C2'-C3'-O3'	5.43	117.64	109.50
27	YD	65	ILE	N-CA-C	5.42	115.96	110.05
28	RE	174	ASP	N-CA-C	5.42	115.40	108.24
10	QJ	53	PRO	CA-C-N	5.42	130.64	123.00
10	QJ	53	PRO	C-N-CA	5.42	130.64	123.00
39	YT	16	ARG	N-CA-C	-5.41	102.19	110.14
12	QL	127	GLU	N-CA-C	-5.40	105.92	113.37
29	YF	169	ASN	N-CA-C	5.39	118.96	112.38
36	RQ	26	TYR	N-CA-C	5.39	118.93	111.71
29	YF	78	ILE	CB-CA-C	-5.38	104.99	112.04
10	XJ	52	GLY	CA-C-N	-5.37	113.09	119.05
10	XJ	52	GLY	C-N-CA	-5.37	113.09	119.05
28	RE	2	LYS	N-CA-C	5.36	118.57	110.64
28	YE	2	LYS	N-CA-C	5.35	118.55	110.64
29	YF	10	PRO	N-CA-C	5.34	118.98	110.21
31	YH	38	SER	CA-C-N	5.34	125.01	119.56
31	YH	38	SER	C-N-CA	5.34	125.01	119.56
29	YF	199	TRP	N-CA-C	-5.34	105.40	112.34
34	RO	32	TYR	N-CA-C	5.34	117.30	109.24
25	RA	271(B)	G	P-O3'-C3'	5.33	126.09	119.70
31	RH	28	GLY	CA-C-N	5.32	126.49	119.84
31	RH	28	GLY	C-N-CA	5.32	126.49	119.84
20	QT	33	ILE	CB-CA-C	-5.31	105.08	111.88
1	XA	812	C	P-O3'-C3'	5.30	128.15	120.20
27	YD	240	ALA	CA-C-N	-5.29	113.22	119.84
27	YD	240	ALA	C-N-CA	-5.29	113.22	119.84
31	YH	28	GLY	CA-C-N	5.28	126.44	119.84
31	YH	28	GLY	C-N-CA	5.28	126.44	119.84
48	Y2	39	ALA	N-CA-C	-5.27	106.80	113.18
25	RA	1130	U	C2'-C3'-O3'	5.27	117.41	109.50
19	XS	41	VAL	C-N-CD	-5.27	109.01	120.60
37	YR	58	GLY	N-CA-C	5.27	122.36	115.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
45	YZ	82	ARG	CA-C-N	5.26	125.51	119.83
45	YZ	82	ARG	C-N-CA	5.26	125.51	119.83
39	RT	13	ARG	N-CA-C	-5.25	106.46	112.92
28	YE	80	GLU	N-CA-C	-5.25	100.88	109.80
27	YD	133	LEU	N-CA-C	5.23	117.66	111.33
37	YR	43	GLU	N-CA-C	-5.23	107.06	113.50
28	YE	200	GLU	N-CA-C	-5.23	102.00	110.32
19	XS	34	TRP	N-CA-C	-5.22	106.93	113.72
50	R4	12	ALA	N-CA-C	5.22	114.73	107.73
27	RD	157	ARG	N-CA-C	-5.22	106.82	114.39
45	RZ	129	SER	CA-C-N	5.22	126.36	119.84
45	RZ	129	SER	C-N-CA	5.22	126.36	119.84
28	RE	80	GLU	N-CA-C	-5.21	100.94	109.80
28	RE	97	LYS	N-CA-C	-5.21	103.13	110.31
19	XS	36	ARG	N-CA-C	-5.20	106.96	113.72
43	RX	5	TYR	N-CA-C	-5.20	107.11	112.93
28	YE	97	LYS	N-CA-C	-5.20	103.14	110.31
28	RE	200	GLU	N-CA-C	-5.20	102.06	110.32
28	YE	80	GLU	CA-C-N	-5.19	117.66	123.16
28	YE	80	GLU	C-N-CA	-5.19	117.66	123.16
39	YT	13	ARG	N-CA-C	-5.19	99.74	110.80
8	QH	26	VAL	CA-C-N	5.19	126.32	119.84
8	QH	26	VAL	C-N-CA	5.19	126.32	119.84
50	R4	18	CYS	CB-CA-C	-5.18	108.91	116.53
28	YE	46	ALA	N-CA-C	5.18	114.68	107.73
4	XD	31	CYS	CA-C-N	5.18	131.43	121.54
4	XD	31	CYS	C-N-CA	5.18	131.43	121.54
28	RE	46	ALA	N-CA-C	5.17	114.66	107.73
7	XG	27	ILE	N-CA-C	-5.17	105.67	110.53
28	YE	194	GLY	N-CA-C	-5.17	105.93	112.54
2	XB	166	ASP	CA-C-N	5.16	124.88	119.82
2	XB	166	ASP	C-N-CA	5.16	124.88	119.82
28	RE	79	ARG	N-CA-C	5.16	121.79	110.80
5	QE	95	ALA	CA-C-N	5.16	126.29	119.84
5	QE	95	ALA	C-N-CA	5.16	126.29	119.84
28	YE	79	ARG	N-CA-C	5.15	121.78	110.80
48	R2	5	GLU	N-CA-C	-5.15	105.74	112.23
45	RZ	100	VAL	CA-C-N	5.15	126.27	119.84
45	RZ	100	VAL	C-N-CA	5.15	126.27	119.84
44	RY	55	TYR	CA-C-N	5.14	126.27	119.84
44	RY	55	TYR	C-N-CA	5.14	126.27	119.84
28	RE	194	GLY	N-CA-C	-5.13	105.97	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
32	RI	79	ILE	N-CA-C	5.13	113.58	107.73
16	QP	39	TYR	N-CA-C	5.13	116.39	107.93
54	R8	53	PRO	N-CA-C	-5.13	101.90	112.47
36	RQ	90	VAL	N-CA-C	5.13	120.00	109.34
50	R4	38	LYS	N-CA-C	-5.13	105.92	113.72
54	Y8	53	PRO	N-CA-C	-5.13	101.91	112.47
36	YQ	90	VAL	N-CA-C	5.12	120.00	109.34
44	YY	76	CYS	CA-C-N	5.12	126.24	119.84
44	YY	76	CYS	C-N-CA	5.12	126.24	119.84
25	RA	1992	G	C2'-C3'-O3'	5.12	117.18	109.50
39	YT	108	ARG	N-CA-C	-5.12	107.16	113.41
25	YA	1992	G	C2'-C3'-O3'	5.12	117.17	109.50
44	YY	60	PHE	N-CA-C	5.12	116.96	109.24
27	RD	256	GLY	N-CA-C	5.11	119.77	113.79
1	XA	1200	C	C3'-C2'-C1'	5.10	106.60	101.50
47	Y1	70	VAL	CB-CA-C	-5.10	105.44	111.97
52	Y6	28	ARG	N-CA-C	5.09	117.13	109.23
29	YF	80	ALA	CA-C-N	5.09	125.37	119.47
29	YF	80	ALA	C-N-CA	5.09	125.37	119.47
48	Y2	56	GLN	N-CA-C	-5.09	105.73	111.28
31	YH	92	ILE	N-CA-C	5.09	119.92	109.34
51	R5	41	PRO	CA-C-N	5.08	126.19	119.84
51	R5	41	PRO	C-N-CA	5.08	126.19	119.84
13	XM	117	VAL	N-CA-C	5.07	117.03	108.81
35	YP	64	LYS	N-CA-C	-5.07	107.14	113.38
28	RE	80	GLU	CA-C-N	-5.06	117.79	123.16
28	RE	80	GLU	C-N-CA	-5.06	117.79	123.16
48	R2	46	GLN	N-CA-C	5.06	119.46	113.28
53	Y7	30	VAL	CB-CA-C	-5.06	105.40	111.88
27	YD	243	GLY	N-CA-C	-5.04	107.84	114.95
29	YF	205	ARG	N-CA-C	-5.04	105.47	110.97
10	XJ	58	ASP	N-CA-C	-5.04	102.18	109.59
20	XT	9	ASN	N-CA-C	5.04	117.94	110.48
27	RD	8	PRO	CA-C-O	-5.04	117.58	121.31
27	RD	196	VAL	N-CA-C	5.04	115.26	110.53
45	RZ	5	LEU	O-C-N	5.02	124.54	120.83
5	XE	127	ASN	CA-C-N	5.02	125.04	119.32
5	XE	127	ASN	C-N-CA	5.02	125.04	119.32
10	QJ	58	ASP	N-CA-C	-5.02	102.22	109.59
27	YD	7	LYS	N-CA-C	-5.00	102.94	110.24

There are no chirality outliers.

There are no planarity outliers.

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	QA	32246	0	16278	532	0
1	XA	32249	0	16279	525	0
2	QB	1924	0	1975	66	0
2	XB	1924	0	1975	90	0
3	QC	1605	0	1668	52	0
3	XC	1605	0	1668	61	0
4	QD	1703	0	1764	60	0
4	XD	1703	0	1764	52	0
5	QE	1155	0	1213	63	0
5	XE	1155	0	1213	59	0
6	QF	843	0	857	18	0
6	XF	843	0	857	22	0
7	QG	1257	0	1296	35	0
7	XG	1257	0	1296	21	0
8	QH	1116	0	1177	42	0
8	XH	1116	0	1177	32	0
9	QI	1010	0	1037	32	0
9	XI	1010	0	1037	49	0
10	QJ	801	0	849	67	0
10	XJ	801	0	849	60	0
11	QK	885	0	904	28	0
11	XK	885	0	904	28	0
12	QL	975	0	1062	103	0
12	XL	975	0	1062	100	0
13	QM	964	0	1034	55	0
13	XM	964	0	1034	45	0
14	QN	492	0	529	22	0
14	XN	492	0	529	22	0
15	QO	734	0	771	20	0
15	XO	734	0	771	17	0
16	QP	705	0	725	14	0
16	XP	705	0	725	23	0
17	QQ	834	0	904	26	0
17	XQ	834	0	904	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	QR	574	0	644	11	0
18	XR	574	0	644	22	0
19	QS	674	0	699	70	0
19	XS	674	0	699	43	0
20	QT	763	0	860	34	0
20	XT	763	0	861	52	0
21	QU	217	0	234	10	0
21	XU	217	0	234	5	0
22	QV	1644	0	836	20	0
22	XV	1644	0	836	17	0
23	QX	220	0	109	1	0
23	XX	220	0	109	3	0
24	QY	321	0	162	2	0
24	XY	344	0	173	8	0
25	RA	62071	0	31291	909	0
25	YA	62071	0	31291	927	0
26	RB	2573	0	1306	41	0
26	YB	2573	0	1306	46	0
27	RD	2115	0	2195	101	0
27	YD	2115	0	2195	336	0
28	RE	1568	0	1634	305	0
28	YE	1568	0	1634	305	0
29	RF	1585	0	1632	55	0
29	YF	1585	0	1632	200	0
30	RG	1474	0	1535	111	0
30	YG	1474	0	1535	66	0
31	RH	1307	0	1382	233	0
31	YH	1307	0	1382	235	0
32	RI	1136	0	1223	74	0
32	YI	1136	0	1223	48	0
33	RN	1104	0	1180	47	0
33	YN	1104	0	1180	57	0
34	RO	933	0	996	26	0
34	YO	933	0	996	27	0
35	RP	1145	0	1228	73	0
35	YP	1145	0	1228	84	0
36	RQ	1122	0	1179	175	0
36	YQ	1122	0	1179	186	0
37	RR	968	0	1033	45	0
37	YR	968	0	1033	35	0
38	RS	882	0	943	46	0
38	YS	882	0	943	169	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
39	RT	1141	0	1202	66	0
39	YT	1141	0	1202	59	0
40	RU	964	0	1022	30	0
40	YU	964	0	1022	56	0
41	RV	779	0	852	19	0
41	YV	779	0	852	50	0
42	RW	900	0	964	24	0
42	YW	900	0	964	25	0
43	RX	725	0	778	29	0
43	YX	725	0	778	21	0
44	RY	785	0	878	53	0
44	YY	785	0	878	42	0
45	RZ	1461	0	1493	43	0
45	YZ	1461	0	1493	63	0
46	R0	648	0	672	30	0
46	Y0	648	0	672	40	0
47	R1	763	0	848	25	0
47	Y1	763	0	848	32	0
48	R2	581	0	629	16	0
48	Y2	581	0	629	79	0
49	R3	469	0	518	8	0
49	Y3	469	0	518	13	0
50	R4	581	0	575	207	0
50	Y4	581	0	577	65	0
51	R5	459	0	480	93	0
51	Y5	454	0	475	39	0
52	R6	424	0	450	32	0
52	Y6	424	0	450	33	0
53	R7	430	0	480	20	0
53	Y7	430	0	480	20	0
54	R8	517	0	582	124	0
54	Y8	517	0	582	127	0
55	R9	307	0	338	11	0
55	Y9	307	0	338	12	0
56	Z5	74	0	51	8	0
56	Z6	74	0	51	11	0
57	QA	95	0	0	0	0
57	QE	1	0	0	0	0
57	QF	1	0	0	0	0
57	QM	1	0	0	0	0
57	QV	3	0	0	0	0
57	QX	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
57	QY	1	0	0	0	0
57	R0	1	0	0	0	0
57	R1	1	0	0	0	0
57	R5	1	0	0	0	0
57	R8	1	0	0	0	0
57	RA	307	0	0	0	0
57	RB	5	0	0	0	0
57	RD	1	0	0	0	0
57	RE	2	0	0	0	0
57	RF	1	0	0	0	0
57	RR	1	0	0	0	0
57	XA	114	0	0	0	0
57	XF	1	0	0	0	0
57	XV	3	0	0	0	0
57	XX	2	0	0	0	0
57	Y0	3	0	0	0	0
57	Y5	1	0	0	0	0
57	Y7	1	0	0	0	0
57	YA	337	0	0	0	0
57	YB	3	0	0	0	0
57	YD	1	0	0	0	0
57	YE	2	0	0	0	0
57	YP	2	0	0	0	0
57	YQ	1	0	0	0	0
58	QA	42	0	45	3	0
58	XA	42	0	45	1	0
59	QD	1	0	0	0	0
59	QN	1	0	0	0	0
59	XD	1	0	0	0	0
59	XN	1	0	0	0	0
All	All	292311	0	198412	8070	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 17.

All (8070) 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)
31:RH:127:GLU:CG	31:RH:128:PRO:HD3	1.35	1.53
31:YH:127:GLU:CG	31:YH:128:PRO:HD3	1.36	1.52
30:RG:3:LEU:HD11	50:R4:25:TYR:CE1	1.65	1.29

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:Y5:49:CYS:SG	51:Y5:60:VAL:HG12	1.75	1.26
32:RI:144:VAL:O	32:RI:145:VAL:HG12	1.31	1.24
36:RQ:59:ARG:O	36:RQ:60:ARG:HD2	1.38	1.22
31:YH:127:GLU:CB	31:YH:128:PRO:HD3	1.69	1.21
31:RH:127:GLU:HG2	31:RH:128:PRO:CD	1.69	1.21
31:YH:127:GLU:HG2	31:YH:128:PRO:CD	1.69	1.21
28:RE:14:ILE:HD11	39:RT:14:TYR:OH	1.35	1.19
31:RH:127:GLU:CB	31:RH:128:PRO:HD3	1.69	1.19
32:RI:53:ALA:O	32:RI:57:ARG:HG2	1.41	1.18
10:QJ:50:ILE:HD11	10:QJ:57:LYS:HD3	1.23	1.16
35:RP:61:ARG:HE	54:R8:13:ARG:HD2	1.07	1.16
31:RH:132:ARG:HH11	31:RH:132:ARG:HB2	1.10	1.15
30:RG:112:PRO:HB3	50:R4:37:SER:HB2	1.27	1.15
10:QJ:50:ILE:HD12	10:QJ:57:LYS:HG2	1.28	1.14
25:YA:242:G:H5"	54:Y8:62:LEU:HD22	1.29	1.13
25:YA:67:U:N3	25:YA:74:A:H2	1.48	1.11
28:YE:179:GLU:HB3	28:YE:181:LEU:HD23	1.32	1.11
31:YH:132:ARG:HB2	31:YH:132:ARG:HH11	1.10	1.11
27:YD:44:ASN:HB2	27:YD:48:ARG:O	1.51	1.11
28:RE:179:GLU:HB3	28:RE:181:LEU:HD23	1.32	1.10
31:RH:86:GLU:HG3	31:RH:165:ALA:H	1.06	1.10
29:YF:101:LEU:HD12	29:YF:102:PRO:HD2	1.21	1.10
45:YZ:182:LYS:HG3	45:YZ:183:LEU:HA	1.21	1.10
28:YE:50:GLY:HA2	28:YE:77:ILE:HA	1.31	1.10
36:YQ:81:VAL:HG23	46:Y0:7:LEU:HD21	1.34	1.10
25:RA:747:U:N3	51:R5:2:ALA:N	2.01	1.09
33:YN:63:THR:HG23	33:YN:66:LYS:HZ2	1.17	1.09
25:YA:593:G:O2'	54:Y8:61:LEU:HD13	1.51	1.09
31:YH:152:ARG:HG3	31:YH:153:LYS:HE2	1.33	1.08
31:YH:86:GLU:HG3	31:YH:165:ALA:H	1.05	1.08
27:YD:131:LEU:HB2	27:YD:136:ILE:HD11	1.35	1.08
31:RH:152:ARG:HG3	31:RH:153:LYS:HE2	1.33	1.07
25:RA:747:U:C4	51:R5:2:ALA:N	2.23	1.07
28:RE:50:GLY:HA2	28:RE:77:ILE:HA	1.31	1.07
28:RE:21:VAL:HB	28:RE:22:PRO:HB3	1.37	1.06
50:R4:71:ARG:HH11	50:R4:71:ARG:HG3	1.13	1.06
28:RE:63:LEU:HD12	28:RE:64:LYS:H	1.18	1.06
54:Y8:52:LYS:H	54:Y8:53:PRO:CD	1.69	1.06
31:YH:153:LYS:HB3	31:YH:154:PRO:HD2	1.07	1.06
28:YE:63:LEU:HD12	28:YE:64:LYS:H	1.18	1.05
28:YE:21:VAL:HB	28:YE:22:PRO:HB3	1.37	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:882:G:H1	25:RA:894:C:N4	1.54	1.05
25:YA:242:G:C5'	54:Y8:62:LEU:HD22	1.86	1.05
25:RA:2015:A:H1'	51:R5:2:ALA:HA	1.34	1.05
31:RH:127:GLU:CB	31:RH:128:PRO:CD	2.35	1.04
36:RQ:59:ARG:O	36:RQ:60:ARG:CD	2.05	1.04
1:XA:838:G:H1	1:XA:848:C:N4	1.54	1.04
31:YH:127:GLU:CG	31:YH:128:PRO:CD	2.31	1.04
53:Y7:9:ARG:NH2	53:Y7:48:LYS:HB2	1.72	1.04
25:YA:1826:G:H4'	27:YD:242:ARG:HH21	1.22	1.04
25:YA:674:G:H1'	29:YF:74:ARG:HD3	1.39	1.04
35:RP:61:ARG:NE	54:R8:13:ARG:HD2	1.71	1.03
36:RQ:81:VAL:O	36:RQ:82:ARG:CD	2.06	1.03
25:YA:2701:C:H3'	25:YA:2702:U:H5"	1.40	1.03
32:RI:53:ALA:O	32:RI:57:ARG:CG	2.07	1.03
38:YS:83:LYS:O	38:YS:109:GLY:HA3	1.57	1.03
31:YH:127:GLU:CB	31:YH:128:PRO:CD	2.35	1.03
38:YS:106:ARG:HA	38:YS:110:LEU:HD11	1.40	1.03
54:R8:52:LYS:H	54:R8:53:PRO:CD	1.69	1.03
10:XJ:50:ILE:CD1	10:XJ:57:LYS:HG2	1.89	1.03
27:YD:35:LYS:HG2	27:YD:64:ILE:N	1.72	1.03
36:RQ:80:GLU:O	36:RQ:81:VAL:HG13	1.59	1.02
36:YQ:81:VAL:O	36:YQ:82:ARG:CD	2.06	1.02
25:RA:1064:C:H42	25:RA:1074:G:H1	1.06	1.02
36:YQ:65:PHE:O	36:YQ:66:ILE:HG12	1.59	1.02
53:R7:9:ARG:HH21	53:R7:48:LYS:HD2	1.18	1.02
29:YF:67:GLN:O	29:YF:68:LYS:HB2	1.56	1.02
36:YQ:12:GLN:HG2	36:YQ:73:PRO:HD2	1.42	1.02
36:RQ:12:GLN:HG2	36:RQ:73:PRO:HD2	1.42	1.02
29:YF:46:ARG:HH11	29:YF:46:ARG:HG2	1.20	1.02
4:QD:9:CYS:SG	4:QD:31:CYS:O	2.18	1.02
31:RH:153:LYS:HB3	31:RH:154:PRO:HD2	1.06	1.02
25:YA:67:U:N3	25:YA:74:A:C2	2.25	1.02
25:RA:2729:G:H1'	28:RE:187:ALA:HB2	1.39	1.01
36:YQ:79:LEU:CD1	46:Y0:5:LYS:HD3	1.90	1.01
36:YQ:79:LEU:HD12	46:Y0:5:LYS:CD	1.90	1.01
5:QE:71:LEU:CD1	5:QE:115:VAL:H	1.72	1.01
25:RA:155:C:H42	25:RA:171:G:H1	1.05	1.01
1:XA:954:G:H21	1:XA:1227:A:H62	1.06	1.01
25:YA:2810:A:O3'	28:YE:61:ARG:HG3	1.60	1.01
29:YF:67:GLN:O	29:YF:67:GLN:HG3	1.58	1.01
53:Y7:9:ARG:HH21	53:Y7:48:LYS:HD2	1.21	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2308:G:H1	25:RA:2311:A:H2	1.09	1.01
30:RG:3:LEU:CD1	50:R4:25:TYR:CE1	2.44	1.00
36:RQ:65:PHE:O	36:RQ:66:ILE:HG12	1.59	1.00
36:RQ:81:VAL:O	36:RQ:82:ARG:NE	1.94	1.00
4:QD:9:CYS:SG	4:QD:22:LYS:HE2	2.02	1.00
25:RA:1359:A:C6	25:RA:1372:U:O4	2.15	1.00
50:R4:56:VAL:HA	50:R4:60:GLN:HB2	1.43	1.00
29:YF:185:ASP:HA	29:YF:188:ARG:HD3	1.41	1.00
36:YQ:79:LEU:HD12	46:Y0:5:LYS:HD3	1.43	1.00
25:RA:2701:C:H3'	25:RA:2702:U:H5"	1.44	1.00
36:YQ:80:GLU:O	36:YQ:81:VAL:HG13	1.59	1.00
31:RH:153:LYS:HB3	31:RH:154:PRO:CD	1.92	0.99
28:RE:201:THR:HG22	28:RE:203:LYS:H	1.26	0.99
31:RH:127:GLU:CG	31:RH:128:PRO:CD	2.31	0.99
38:YS:26:LEU:HD12	38:YS:39:ILE:HD11	1.40	0.99
48:Y2:50:ILE:HD12	48:Y2:51:ARG:N	1.76	0.99
31:RH:77:LYS:HB3	31:RH:77:LYS:HZ3	1.22	0.99
28:YE:201:THR:HG22	28:YE:203:LYS:H	1.26	0.99
31:YH:153:LYS:HB3	31:YH:154:PRO:CD	1.92	0.99
35:RP:62:LEU:CD2	54:R8:25:MET:HB2	1.91	0.99
25:RA:2068:U:H3	25:RA:2430:A:H2	1.09	0.99
30:RG:112:PRO:HB3	50:R4:37:SER:CB	1.92	0.99
30:RG:5:VAL:HG22	50:R4:25:TYR:CD2	1.97	0.99
51:Y5:49:CYS:SG	51:Y5:60:VAL:CG1	2.50	0.99
30:RG:3:LEU:HD11	50:R4:25:TYR:HE1	0.97	0.98
30:RG:179:PRO:HG3	50:R4:38:LYS:HZ2	1.28	0.98
25:YA:2099:U:H3	25:YA:2190:G:H1	1.01	0.98
36:YQ:81:VAL:O	36:YQ:82:ARG:NE	1.94	0.98
32:RI:144:VAL:O	32:RI:145:VAL:CG1	2.12	0.98
36:RQ:79:LEU:O	36:RQ:79:LEU:HD22	1.64	0.98
31:YH:86:GLU:HG3	31:YH:165:ALA:N	1.79	0.98
25:YA:592:G:H21	54:Y8:4:MET:HE2	1.25	0.98
10:QJ:50:ILE:HD11	10:QJ:57:LYS:CD	1.93	0.97
25:RA:1169:G:H1	25:RA:1180:C:N4	1.61	0.97
36:YQ:79:LEU:O	36:YQ:79:LEU:HD13	1.63	0.97
13:QM:3:ARG:HH22	30:RG:113:ARG:HH21	1.11	0.97
28:RE:20:ALA:O	28:RE:21:VAL:HG22	1.64	0.97
48:Y2:50:ILE:HD12	48:Y2:51:ARG:H	1.24	0.97
25:RA:873:G:H1	25:RA:904:C:H42	0.97	0.97
27:YD:44:ASN:HB3	27:YD:49:ILE:HA	1.45	0.97
38:YS:83:LYS:NZ	38:YS:109:GLY:HA2	1.78	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:YQ:79:LEU:O	36:YQ:79:LEU:HD22	1.64	0.97
25:RA:2572:A:C2	28:RE:144:ARG:NH2	2.33	0.97
13:QM:3:ARG:NH2	30:RG:113:ARG:HH21	1.62	0.96
30:RG:3:LEU:HD21	50:R4:25:TYR:CD1	2.00	0.96
36:RQ:79:LEU:O	36:RQ:79:LEU:HD13	1.63	0.96
25:YA:784:A:N7	27:YD:229:VAL:HG21	1.80	0.96
5:XE:41:VAL:CG1	5:XE:113:ALA:HB2	1.94	0.96
25:RA:2107:C:H42	25:RA:2182:G:H1	1.11	0.96
28:YE:20:ALA:O	28:YE:21:VAL:HG22	1.65	0.96
36:RQ:79:LEU:HD12	46:R0:5:LYS:HD3	1.46	0.95
25:YA:330:A:H2	25:YA:1210:A:HO2'	1.07	0.95
29:YF:101:LEU:HD12	29:YF:102:PRO:CD	1.96	0.95
12:QL:6:THR:H	12:QL:9:GLN:HE21	1.15	0.95
20:XT:26:ASN:HB2	20:XT:71:THR:HG23	1.48	0.95
29:YF:103:LYS:HA	29:YF:106:ARG:HG3	1.48	0.95
25:YA:2308:G:H1	25:YA:2311:A:H2	0.98	0.95
31:RH:86:GLU:HG3	31:RH:165:ALA:N	1.79	0.95
27:YD:227:ASN:HB3	27:YD:228:PRO:HD2	1.44	0.94
28:YE:78:LEU:HG	28:YE:79:ARG:HE	1.30	0.94
31:YH:153:LYS:CB	31:YH:154:PRO:HD2	1.98	0.94
48:Y2:13:ALA:HA	48:Y2:16:LEU:HD23	1.48	0.94
50:R4:36:CYS:O	50:R4:39:CYS:HB2	1.67	0.94
10:QJ:50:ILE:CD1	10:QJ:57:LYS:HG2	1.96	0.94
28:YE:14:ILE:HD11	39:YT:14:TYR:OH	1.65	0.94
25:RA:2287:A:H62	25:RA:2344:U:H3	1.10	0.94
36:RQ:80:GLU:OE1	46:R0:7:LEU:HB3	1.67	0.94
28:RE:78:LEU:HG	28:RE:79:ARG:HE	1.31	0.94
25:YA:483:A:H4'	44:YY:49:VAL:HA	1.50	0.94
31:YH:77:LYS:NZ	31:YH:77:LYS:HB3	1.82	0.94
30:RG:65:GLY:O	50:R4:7:PRO:HD2	1.66	0.94
27:YD:28:GLU:HB2	27:YD:29:PRO:CD	1.98	0.94
31:YH:77:LYS:HB3	31:YH:77:LYS:HZ3	1.31	0.94
5:QE:69:VAL:HG12	5:QE:71:LEU:HD23	1.48	0.93
19:QS:68:GLY:CA	50:R4:68:ARG:HG2	1.99	0.93
2:QB:185:ILE:HG22	2:QB:199:TYR:HB2	1.48	0.93
51:R5:56:LYS:H	51:R5:56:LYS:HD2	1.31	0.93
35:RP:62:LEU:HD21	54:R8:25:MET:HB2	1.51	0.93
36:RQ:59:ARG:O	36:RQ:60:ARG:CG	2.17	0.93
31:YH:127:GLU:HB3	31:YH:128:PRO:CD	1.99	0.93
40:YU:90:VAL:HG22	41:YV:39:LEU:HB3	1.51	0.93
25:YA:1803:A:H4'	27:YD:259:THR:CG2	1.99	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:RQ:34:LEU:HD11	36:RQ:129:THR:HB	1.50	0.93
1:XA:148:G:H1	1:XA:174:C:H42	0.99	0.93
1:XA:80:G:H1	1:XA:89:U:H3	1.04	0.93
27:YD:108:PRO:HB3	27:YD:143:HIS:HE1	1.32	0.93
1:QA:1002:G:H1	1:QA:1038:C:H42	0.99	0.93
31:RH:127:GLU:HB3	31:RH:128:PRO:CD	1.99	0.93
27:YD:108:PRO:HB3	27:YD:143:HIS:CE1	2.05	0.93
51:R5:58:LEU:HD13	51:R5:60:VAL:HG12	1.48	0.92
1:XA:838:G:H1	1:XA:848:C:H42	1.06	0.92
38:YS:59:LYS:HG2	38:YS:60:GLY:H	1.31	0.92
25:YA:1169:G:H1	25:YA:1180:C:N4	1.68	0.92
28:YE:14:ILE:HG12	28:YE:15:PHE:H	1.33	0.92
25:YA:1169:G:H1	25:YA:1180:C:H42	0.98	0.92
25:RA:2245:U:H5'	25:RA:2246:G:H5'	1.52	0.92
31:RH:153:LYS:CB	31:RH:154:PRO:HD2	1.97	0.92
25:RA:2810:A:O3'	28:RE:61:ARG:HG3	1.69	0.91
30:RG:113:ARG:HG2	50:R4:34:GLU:CD	1.95	0.91
25:YA:631:A:OP2	54:Y8:46:ARG:NH2	2.01	0.91
38:YS:67:ARG:NH1	38:YS:67:ARG:HB2	1.85	0.91
36:YQ:34:LEU:HD11	36:YQ:129:THR:HB	1.50	0.91
31:RH:77:LYS:HB3	31:RH:77:LYS:NZ	1.82	0.91
27:YD:10:THR:HG23	27:YD:13:ARG:HB3	1.51	0.91
25:RA:301:G:H1	25:RA:316:C:H42	1.19	0.91
10:XJ:50:ILE:HD11	10:XJ:57:LYS:HG2	1.53	0.91
25:YA:1138:G:H21	33:YN:106:MET:HE3	1.35	0.91
28:RE:14:ILE:HG12	28:RE:15:PHE:H	1.33	0.91
33:YN:7:LYS:H	33:YN:7:LYS:HD2	1.36	0.90
25:RA:2808:U:H3	25:RA:2892:A:H62	1.18	0.90
25:YA:1113:U:OP1	31:YH:2:SER:N	2.04	0.90
25:YA:2308:G:N1	25:YA:2311:A:C2	2.39	0.90
48:Y2:65:ASN:HB3	48:Y2:69:ARG:HH12	1.33	0.90
25:RA:1113:U:OP1	31:RH:2:SER:N	2.03	0.90
31:RH:4:ILE:HG13	31:RH:6:ARG:CZ	2.01	0.90
19:QS:5:LEU:HD21	50:R4:67:TYR:CZ	2.06	0.90
25:RA:873:G:H1	25:RA:904:C:N4	1.69	0.90
31:YH:4:ILE:HG13	31:YH:6:ARG:CZ	2.01	0.90
36:RQ:59:ARG:O	36:RQ:60:ARG:HG3	1.72	0.90
31:YH:26:VAL:HG13	31:YH:27:LYS:H	1.35	0.90
20:XT:41:ILE:HG22	20:XT:91:LEU:HD12	1.54	0.90
25:YA:1359:A:N6	25:YA:1372:U:H3	1.67	0.90
30:RG:6:ALA:CB	50:R4:23:GLU:HG2	2.03	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:RG:6:ALA:H	50:R4:23:GLU:CG	1.85	0.89
33:YN:63:THR:CG2	33:YN:66:LYS:HZ2	1.84	0.89
54:Y8:52:LYS:H	54:Y8:53:PRO:HD3	1.35	0.89
30:RG:113:ARG:HG2	50:R4:34:GLU:OE2	1.72	0.89
25:YA:2287:A:H62	25:YA:2344:U:H3	1.17	0.89
27:YD:69:ARG:HH21	27:YD:130:ALA:HB2	1.37	0.89
27:YD:147:LEU:HD13	27:YD:155:LEU:HD11	1.51	0.89
25:RA:1359:A:N1	25:RA:1372:U:O4	2.04	0.89
31:RH:10:PRO:HD2	31:RH:50:VAL:O	1.72	0.89
31:RH:26:VAL:HG13	31:RH:27:LYS:H	1.36	0.89
12:XL:6:THR:H	12:XL:9:GLN:HE21	1.15	0.89
27:YD:44:ASN:HD22	27:YD:44:ASN:H	1.19	0.89
27:YD:183:ARG:HG2	27:YD:183:ARG:HH11	1.34	0.89
36:YQ:81:VAL:CG2	46:Y0:7:LEU:HD21	2.02	0.89
1:QA:589:C:H42	1:QA:650:G:H1	1.21	0.89
51:R5:3:LYS:HA	51:R5:3:LYS:HE3	1.54	0.89
54:Y8:59:LYS:NZ	54:Y8:59:LYS:HB2	1.88	0.89
28:YE:63:LEU:HD12	28:YE:64:LYS:N	1.87	0.89
25:RA:1043:C:H42	25:RA:1112:G:H1	1.19	0.89
25:RA:2014:A:O2'	51:R5:2:ALA:HB2	1.73	0.89
1:QA:1395:C:HO2'	1:QA:1401:G:HO2'	1.19	0.89
54:R8:52:LYS:H	54:R8:53:PRO:HD3	1.35	0.88
31:YH:10:PRO:HD2	31:YH:50:VAL:O	1.72	0.88
44:YY:79:CYS:SG	44:YY:80:GLY:N	2.45	0.88
5:XE:41:VAL:HG13	5:XE:113:ALA:HB2	1.53	0.88
5:QE:101:ILE:HD13	5:QE:101:ILE:H	1.36	0.88
39:RT:26:ASP:HB3	39:RT:92:GLY:H	1.36	0.88
25:RA:155:C:N4	25:RA:171:G:H1	1.72	0.88
26:RB:51:G:N7	38:RS:62:LYS:NZ	2.20	0.88
25:RA:1443:G:H1	25:RA:1548:C:H42	1.19	0.88
19:QS:5:LEU:CD2	50:R4:67:TYR:CZ	2.56	0.88
28:RE:63:LEU:HD12	28:RE:64:LYS:N	1.88	0.88
25:YA:528:A:O2'	25:YA:529:A:H5'	1.74	0.88
25:YA:593:G:O3'	54:Y8:61:LEU:HD22	1.73	0.88
25:YA:242:G:H5"	54:Y8:62:LEU:CD2	2.04	0.88
25:YA:2245:U:H5'	25:YA:2246:G:H5'	1.56	0.88
38:YS:106:ARG:NH1	38:YS:106:ARG:HB2	1.88	0.88
51:R5:40:LYS:HZ1	51:R5:48:GLU:HB2	1.39	0.88
27:YD:27:THR:HG23	27:YD:28:GLU:H	1.38	0.88
25:RA:2308:G:N1	25:RA:2311:A:C2	2.42	0.87
28:RE:77:ILE:HD12	28:RE:78:LEU:N	1.89	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:R8:59:LYS:HB2	54:R8:59:LYS:NZ	1.88	0.87
28:YE:77:ILE:HD12	28:YE:78:LEU:N	1.89	0.87
36:YQ:64:ILE:HA	36:YQ:106:VAL:HG12	1.54	0.87
27:YD:44:ASN:CB	27:YD:49:ILE:HA	2.05	0.87
3:QC:162:GLN:NE2	3:QC:162:GLN:HA	1.88	0.87
36:RQ:64:ILE:HA	36:RQ:106:VAL:HG12	1.54	0.87
27:YD:181:GLU:HA	27:YD:272:ALA:HB3	1.57	0.87
44:YY:76:CYS:HB3	44:YY:96:ILE:HD13	1.57	0.87
25:YA:2056:G:N2	51:Y5:4:HIS:O	2.08	0.86
51:Y5:40:LYS:HG2	51:Y5:47:PRO:HD2	1.56	0.86
30:RG:179:PRO:HG3	50:R4:38:LYS:NZ	1.89	0.86
25:YA:1338:G:N7	43:YX:62:LYS:NZ	2.23	0.86
29:YF:7:TYR:HB3	29:YF:21:ALA:HB1	1.53	0.86
10:QJ:55:LYS:HE3	10:QJ:56:HIS:NE2	1.90	0.86
35:RP:58:THR:O	35:RP:61:ARG:NH2	2.08	0.86
7:XG:89:MET:HE1	7:XG:156:TRP:H	1.41	0.86
25:YA:847:U:O4	25:YA:933:A:N1	2.08	0.86
25:YA:1359:A:N6	25:YA:1372:U:N3	2.22	0.86
31:YH:127:GLU:HG2	31:YH:128:PRO:HD3	0.86	0.86
1:QA:1002:G:H1	1:QA:1038:C:N4	1.73	0.86
28:RE:14:ILE:HD11	39:RT:14:TYR:HH	1.41	0.86
27:YD:35:LYS:HG2	27:YD:64:ILE:H	1.40	0.86
31:RH:8:PRO:C	31:RH:9:ILE:HG12	2.00	0.86
1:XA:148:G:H1	1:XA:174:C:N4	1.72	0.86
27:RD:43:ARG:NH1	27:RD:44:ASN:OD1	2.08	0.86
25:RA:242:G:C5'	54:R8:62:LEU:HD22	2.06	0.85
28:RE:61:ARG:O	28:RE:63:LEU:HG	1.77	0.85
36:RQ:75:THR:HA	36:RQ:88:GLY:O	1.76	0.85
45:RZ:110:GLY:HA2	45:RZ:111:VAL:O	1.75	0.85
30:YG:6:ALA:H	50:Y4:23:GLU:HG2	1.42	0.85
36:YQ:75:THR:HA	36:YQ:88:GLY:O	1.76	0.85
36:YQ:81:VAL:HG23	46:Y0:7:LEU:CD2	2.06	0.85
45:YZ:151:HIS:HB3	45:YZ:170:THR:HA	1.59	0.85
25:RA:2277:G:OP1	36:RQ:85:LYS:HB2	1.76	0.85
1:XA:201:C:H42	1:XA:216:G:H1	1.24	0.85
1:XA:1305:G:N2	1:XA:1331:G:H2'	1.91	0.85
10:XJ:55:LYS:HE3	10:XJ:56:HIS:NE2	1.90	0.85
38:YS:106:ARG:HB2	38:YS:106:ARG:HH11	1.40	0.85
25:RA:1247:A:OP2	35:RP:15:ARG:NH2	2.09	0.85
28:RE:27:LEU:HD21	39:RT:1:MET:HE1	1.57	0.85
25:YA:49:A:N7	25:YA:120:U:C5	2.44	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:RR:33:ARG:NH2	51:R5:55:ARG:HG2	1.91	0.85
51:R5:39:MET:O	51:R5:40:LYS:HG3	1.77	0.85
25:YA:1359:A:H3'	25:YA:1360:A:H5'	1.59	0.85
31:YH:89:ILE:HD11	31:YH:129:THR:HB	1.59	0.85
5:QE:50:GLU:HB3	5:QE:53:LEU:HD13	1.59	0.85
25:RA:2572:A:N3	28:RE:144:ARG:NH2	2.24	0.85
45:RZ:111:VAL:HG22	45:RZ:112:ARG:H	1.39	0.85
22:XV:5:G:H1	22:XV:67:C:H42	1.22	0.85
31:YH:8:PRO:C	31:YH:9:ILE:HG12	2.00	0.85
28:RE:81:ILE:O	28:RE:82:ARG:HB2	1.75	0.85
28:RE:95:ILE:HD12	28:RE:95:ILE:H	1.41	0.85
27:YD:28:GLU:HB2	27:YD:29:PRO:HD2	1.56	0.85
29:YF:82:ILE:HG13	29:YF:82:ILE:O	1.73	0.85
42:YW:18:ARG:HG3	42:YW:76:VAL:HG13	1.58	0.85
54:Y8:59:LYS:NZ	54:Y8:59:LYS:CB	2.40	0.85
31:RH:127:GLU:HG2	31:RH:128:PRO:HD3	0.86	0.85
1:XA:838:G:N2	1:XA:848:C:N3	2.25	0.84
38:YS:83:LYS:HG2	38:YS:109:GLY:CA	2.07	0.84
25:RA:1169:G:H1	25:RA:1180:C:H42	0.85	0.84
28:RE:24:THR:HG21	28:RE:188:VAL:HG11	1.59	0.84
31:RH:89:ILE:HD11	31:RH:129:THR:HB	1.58	0.84
44:RY:79:CYS:SG	44:RY:80:GLY:N	2.45	0.84
28:YE:95:ILE:H	28:YE:95:ILE:HD12	1.41	0.84
31:YH:54:ARG:NH1	31:YH:62:LYS:HG2	1.92	0.84
54:R8:59:LYS:NZ	54:R8:59:LYS:CB	2.40	0.84
25:YA:784:A:C5	27:YD:229:VAL:HG21	2.11	0.84
25:YA:2485:G:OP1	36:YQ:46:GLN:NE2	2.09	0.84
27:YD:17:THR:HG22	27:YD:205:VAL:H	1.41	0.84
38:YS:106:ARG:HA	38:YS:110:LEU:CD1	2.08	0.84
25:RA:2502:G:H5"	25:RA:2503:A:H5"	1.59	0.84
30:RG:112:PRO:CB	50:R4:37:SER:HB2	2.05	0.84
35:YP:59:LEU:O	54:Y8:13:ARG:NH1	2.09	0.84
1:XA:448:A:OP2	1:XA:485:G:N2	2.11	0.84
28:YE:61:ARG:O	28:YE:63:LEU:HG	1.77	0.84
28:YE:81:ILE:O	28:YE:82:ARG:HB2	1.76	0.84
29:YF:53:THR:HG23	29:YF:56:GLU:OE1	1.77	0.84
36:YQ:30:GLY:HA2	36:YQ:107:ALA:HB2	1.60	0.84
25:RA:1899:G:H21	25:RA:1902:C:N4	1.76	0.84
30:YG:98:ARG:HH12	50:Y4:1:MET:HB3	1.41	0.84
39:YT:26:ASP:HB3	39:YT:92:GLY:H	1.42	0.84
19:QS:5:LEU:CD1	50:R4:67:TYR:CE2	2.60	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:RH:54:ARG:NH1	31:RH:62:LYS:HG2	1.92	0.84
35:RP:59:LEU:O	54:R8:13:ARG:NH1	2.10	0.84
35:RP:126:VAL:HG12	35:RP:147:LEU:HD21	1.59	0.84
38:YS:89:ARG:HD2	38:YS:92:TYR:O	1.78	0.84
26:YB:50:G:OP1	38:YS:63:THR:HG23	1.78	0.83
31:RH:105:LEU:HD13	31:RH:105:LEU:H	1.42	0.83
35:RP:61:ARG:HE	54:R8:13:ARG:CD	1.91	0.83
1:XA:1305:G:H22	1:XA:1331:G:H2'	1.41	0.83
2:QB:80:ILE:HD11	2:QB:208:ILE:HG23	1.58	0.83
53:R7:9:ARG:NH2	53:R7:48:LYS:HD2	1.93	0.83
29:YF:32:LEU:HD13	29:YF:105:VAL:HG13	1.59	0.83
25:RA:270(G):C:H42	25:RA:270(S):G:H1	1.21	0.83
30:RG:67:LYS:HG2	50:R4:5:ILE:O	1.78	0.83
31:YH:13:LYS:HE2	31:YH:13:LYS:HA	1.60	0.83
40:YU:88:ILE:HG22	40:YU:90:VAL:HG23	1.59	0.83
25:YA:2747:G:OP1	31:YH:138:LYS:NZ	2.11	0.83
50:Y4:18:CYS:CB	50:Y4:39:CYS:HB3	2.08	0.83
1:QA:671:G:H1	1:QA:735:C:H42	1.26	0.83
51:R5:40:LYS:HD3	51:R5:46:CYS:HB3	1.60	0.83
28:RE:35:GLN:HG2	28:RE:37:ARG:HE	1.44	0.83
28:YE:7:VAL:HG23	28:YE:8:LYS:H	1.44	0.83
25:RA:1138:G:H21	33:RN:106:MET:HE3	1.42	0.83
30:YG:27:ASN:HB3	30:YG:30:GLU:HG3	1.61	0.83
1:QA:49:U:C5	1:QA:365:U:O4	2.32	0.83
1:QA:677:U:H3	1:QA:713:G:H22	1.23	0.83
12:QL:38:THR:HG23	12:QL:39:VAL:HG23	1.60	0.83
25:RA:1310:G:OP2	53:R7:9:ARG:NH1	2.10	0.83
31:RH:153:LYS:HA	31:RH:153:LYS:HZ3	1.43	0.82
36:RQ:30:GLY:HA2	36:RQ:107:ALA:HB2	1.60	0.82
1:XA:664:G:H22	1:XA:741:G:H1	1.26	0.82
1:QA:1453:G:H8	20:QT:39:LYS:HZ1	1.25	0.82
25:RA:2056:G:N2	51:R5:4:HIS:O	2.12	0.82
27:YD:25:THR:CG2	27:YD:82:ILE:H	1.93	0.82
1:XA:339:C:OP2	34:YO:97:ARG:NH1	2.12	0.82
28:YE:15:PHE:CE1	28:YE:20:ALA:HB2	2.14	0.82
31:YH:153:LYS:HG2	31:YH:162:ILE:HG13	1.61	0.82
48:Y2:16:LEU:HG	48:Y2:16:LEU:O	1.78	0.82
25:RA:819:A:OP2	25:RA:1187:G:N2	2.10	0.82
25:RA:1262:A:N3	51:R5:10:LYS:HE3	1.94	0.82
28:RE:15:PHE:CE1	28:RE:20:ALA:HB2	2.14	0.82
31:YH:105:LEU:HD13	31:YH:105:LEU:H	1.42	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:XL:38:THR:HG23	12:XL:39:VAL:HG23	1.60	0.82
10:QJ:50:ILE:HD12	10:QJ:57:LYS:CG	2.09	0.82
45:YZ:182:LYS:CG	45:YZ:183:LEU:HA	2.08	0.82
25:RA:847:U:C4	25:RA:933:A:N1	2.48	0.82
28:RE:7:VAL:HG23	28:RE:8:LYS:H	1.44	0.82
44:RY:29:GLU:HB3	44:RY:38:ILE:HG12	1.62	0.82
54:R8:59:LYS:HB2	54:R8:59:LYS:HZ2	1.44	0.82
38:YS:19:LYS:O	38:YS:20:ARG:HB3	1.80	0.82
47:Y1:7:ILE:HD12	47:Y1:62:VAL:HG11	1.62	0.82
29:YF:192:LEU:HD21	29:YF:194:MET:HE3	1.60	0.82
25:YA:498:G:N3	44:YY:47:LYS:NZ	2.28	0.81
27:YD:35:LYS:NZ	27:YD:104:TYR:HB2	1.93	0.81
12:QL:86:ARG:HB2	12:QL:101:VAL:HG22	1.62	0.81
51:Y5:16:ARG:NH1	51:Y5:17:ASP:OD1	2.13	0.81
1:QA:346:G:OP1	39:RT:41:ARG:NH2	2.13	0.81
19:QS:68:GLY:HA3	50:R4:68:ARG:HB2	1.62	0.81
31:RH:132:ARG:HB2	31:RH:132:ARG:NH1	1.95	0.81
50:R4:33:VAL:HG12	50:R4:34:GLU:H	1.44	0.81
1:XA:954:G:H4'	13:XM:121:LYS:HG3	1.61	0.81
38:YS:88:ASP:O	38:YS:89:ARG:HB3	1.78	0.81
25:RA:1093:G:OP1	31:RH:170:ARG:HD2	1.80	0.81
31:RH:153:LYS:HG2	31:RH:162:ILE:HG13	1.61	0.81
36:RQ:90:VAL:HG13	36:RQ:91:GLU:N	1.95	0.81
25:YA:265:A:N6	25:YA:427:U:O2'	2.13	0.81
28:YE:24:THR:HG21	28:YE:188:VAL:HG11	1.59	0.81
36:YQ:90:VAL:HG13	36:YQ:91:GLU:N	1.95	0.81
31:RH:152:ARG:O	31:RH:153:LYS:HB2	1.80	0.81
25:RA:2392:A:H1'	35:RP:60:MET:HE3	1.63	0.81
31:YH:10:PRO:O	31:YH:11:VAL:HG13	1.80	0.81
28:RE:3:GLY:O	28:RE:4:ILE:HB	1.81	0.81
54:R8:52:LYS:N	54:R8:53:PRO:CD	2.43	0.81
25:YA:2729:G:H1'	28:YE:187:ALA:HB2	1.61	0.81
28:YE:50:GLY:CA	28:YE:77:ILE:HA	2.09	0.81
48:Y2:43:GLN:O	48:Y2:44:LEU:HG	1.81	0.81
2:XB:69:LEU:HB3	2:XB:162:ILE:HG22	1.62	0.81
38:YS:36:TYR:HD2	38:YS:52:SER:HB3	1.46	0.81
25:RA:242:G:H5"	54:R8:62:LEU:HD22	1.62	0.81
31:RH:13:LYS:HE2	31:RH:13:LYS:HA	1.60	0.81
5:XE:35:GLY:HA3	5:XE:112:LEU:O	1.80	0.81
28:YE:116:VAL:HG21	28:YE:122:PHE:CD2	2.16	0.81
1:QA:1346:A:H5"	9:QI:120:ARG:HH12	1.46	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:QS:5:LEU:HD22	50:R4:67:TYR:OH	1.81	0.81
25:RA:1064:C:N4	25:RA:1074:G:H1	1.76	0.81
25:RA:2807:G:N1	25:RA:2893:G:O6	2.12	0.81
25:RA:2128:C:H42	25:RA:2160:G:H1	1.29	0.80
25:RA:2392:A:H8	35:RP:60:MET:HG2	1.45	0.80
28:RE:116:VAL:HG21	28:RE:122:PHE:CD2	2.16	0.80
13:QM:14:ARG:H	13:QM:44:ARG:HD3	1.45	0.80
28:RE:52:LEU:HB2	28:RE:75:VAL:HG23	1.62	0.80
29:YF:155:LEU:HD13	29:YF:174:VAL:HG13	1.62	0.80
39:RT:24:PRO:HA	39:RT:49:VAL:HG13	1.61	0.80
51:R5:4:HIS:HB3	51:R5:5:PRO:CD	2.11	0.80
25:YA:819:A:OP2	25:YA:1187:G:N2	2.12	0.80
27:YD:27:THR:HG23	27:YD:28:GLU:N	1.96	0.80
1:QA:189:U:O2'	1:QA:190:G:OP1	1.98	0.80
25:RA:882:G:N2	25:RA:894:C:N3	2.27	0.80
28:RE:50:GLY:CA	28:RE:77:ILE:HA	2.10	0.80
33:RN:42:TRP:O	40:RU:64:ARG:NH2	2.15	0.80
1:XA:619:U:H3	4:XD:135:LEU:HD23	1.47	0.80
28:YE:52:LEU:HB2	28:YE:75:VAL:HG23	1.61	0.80
36:YQ:80:GLU:O	36:YQ:81:VAL:CG1	2.30	0.80
38:YS:106:ARG:CA	38:YS:110:LEU:HD21	2.10	0.80
28:RE:201:THR:CG2	28:RE:203:LYS:HB3	2.12	0.80
30:YG:179:PRO:HG3	50:Y4:38:LYS:NZ	1.97	0.80
54:Y8:52:LYS:N	54:Y8:53:PRO:CD	2.43	0.80
25:RA:1045:A:C8	25:RA:1111:A:N6	2.50	0.80
32:RI:54:GLN:O	32:RI:58:LEU:CB	2.28	0.80
28:YE:3:GLY:O	28:YE:4:ILE:HB	1.81	0.80
28:YE:201:THR:CG2	28:YE:203:LYS:HB3	2.11	0.80
29:YF:198:ALA:HA	29:YF:201:VAL:HG12	1.62	0.80
31:YH:152:ARG:O	31:YH:153:LYS:HB2	1.80	0.80
44:YY:76:CYS:SG	44:YY:77:PRO:HD2	2.22	0.80
25:RA:2667:C:H1'	31:RH:109:PHE:HD2	1.46	0.80
31:RH:150:ALA:C	31:RH:152:ARG:H	1.88	0.80
51:R5:20:ARG:HA	51:R5:23:HIS:HD1	1.47	0.80
26:YB:75:G:H5"	45:YZ:36:LYS:HE2	1.64	0.80
27:YD:121:PRO:HB3	27:YD:135:PHE:HE2	1.46	0.80
28:YE:35:GLN:HG2	28:YE:37:ARG:HE	1.44	0.80
13:QM:3:ARG:HA	13:QM:9:ILE:HG21	1.62	0.80
30:YG:98:ARG:NH1	50:Y4:1:MET:SD	2.55	0.80
13:XM:65:LYS:HD3	13:XM:69:GLU:HG3	1.63	0.80
48:R2:46:GLN:O	48:R2:47:ASN:HB2	1.79	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:XD:31:CYS:O	4:XD:31:CYS:SG	2.40	0.79
27:YD:68:LYS:HB2	27:YD:70:TRP:CH2	2.17	0.79
1:XA:1286:A:H5"	21:XU:26:LYS:HD2	1.62	0.79
35:YP:62:LEU:HD21	54:Y8:25:MET:HB2	1.65	0.79
25:RA:530:G:O2'	25:RA:532:A:N7	2.16	0.79
12:XL:86:ARG:HB2	12:XL:101:VAL:HG22	1.62	0.79
31:YH:126:PRO:CG	31:YH:127:GLU:H	1.95	0.79
33:YN:63:THR:HG23	33:YN:66:LYS:NZ	1.97	0.79
11:QK:21:ILE:HB	11:QK:84:VAL:HG12	1.65	0.79
25:RA:847:U:O4	25:RA:933:A:N1	2.16	0.79
28:RE:137:HIS:HB3	28:RE:138:PRO:HD2	1.65	0.79
36:RQ:81:VAL:O	36:RQ:82:ARG:CG	2.31	0.79
3:XC:32:LEU:HD13	3:XC:59:ARG:HD3	1.64	0.79
27:YD:34:VAL:O	27:YD:34:VAL:HG13	1.81	0.79
31:YH:150:ALA:C	31:YH:152:ARG:H	1.88	0.79
31:YH:169:VAL:HG22	31:YH:170:ARG:H	1.48	0.79
38:YS:111:GLU:OE1	38:YS:111:GLU:HA	1.80	0.79
43:YX:67:GLY:O	43:YX:69:TYR:N	2.15	0.79
25:YA:592:G:N2	54:Y8:4:MET:HE2	1.97	0.79
25:YA:674:G:C1'	29:YF:74:ARG:HD3	2.12	0.79
31:YH:26:VAL:HG13	31:YH:27:LYS:N	1.96	0.79
36:YQ:81:VAL:O	36:YQ:82:ARG:CG	2.31	0.79
1:QA:1135:U:H4'	1:QA:1136:U:H5	1.47	0.79
31:RH:10:PRO:O	31:RH:11:VAL:HG13	1.80	0.79
37:RR:104:ARG:HD2	37:RR:111:LEU:HD21	1.63	0.79
28:YE:24:THR:HG21	28:YE:188:VAL:CG1	2.13	0.79
15:QO:26:GLU:OE2	15:QO:77:ARG:NH1	2.16	0.79
31:RH:126:PRO:CG	31:RH:127:GLU:H	1.96	0.79
25:YA:270(T):G:H5"	47:Y1:97:LEU:HD22	1.63	0.79
27:YD:17:THR:CG2	27:YD:205:VAL:H	1.96	0.79
1:QA:664:G:H22	1:QA:741:G:H1	1.29	0.79
12:QL:10:LEU:HB3	17:QQ:32:TYR:CE2	2.18	0.79
25:RA:270(R):G:N3	47:R1:78:LYS:NZ	2.31	0.79
28:RE:111:ARG:HE	28:RE:160:TYR:HE1	1.31	0.79
30:RG:6:ALA:H	50:R4:23:GLU:HG2	1.46	0.79
31:RH:86:GLU:CG	31:RH:165:ALA:H	1.94	0.79
12:XL:6:THR:N	12:XL:9:GLN:HE21	1.80	0.79
31:YH:153:LYS:CG	31:YH:162:ILE:H	1.96	0.79
35:YP:58:THR:O	35:YP:61:ARG:NE	2.15	0.79
51:Y5:49:CYS:SG	51:Y5:60:VAL:N	2.52	0.79
2:XB:77:ALA:HB2	2:XB:211:ILE:HD13	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:530:G:C2	25:YA:2022:U:OP1	2.36	0.78
25:YA:1803:A:H4'	27:YD:259:THR:HG21	1.64	0.78
12:QL:6:THR:N	12:QL:9:GLN:HE21	1.80	0.78
25:RA:1019:U:H3	25:RA:1142(A):A:H62	1.30	0.78
36:RQ:119:ARG:HG2	36:RQ:119:ARG:HH11	1.48	0.78
1:XA:1124:G:H3'	1:XA:1145:C:H41	1.48	0.78
25:YA:592:G:H21	54:Y8:4:MET:CE	1.96	0.78
45:YZ:181:GLU:HG2	45:YZ:183:LEU:HG	1.65	0.78
20:QT:100:ILE:HG13	20:QT:102:GLY:H	1.48	0.78
25:RA:1169:G:N2	25:RA:1180:C:N3	2.29	0.78
36:RQ:80:GLU:O	36:RQ:81:VAL:CG1	2.30	0.78
1:XA:542:G:OP1	4:XD:10:ARG:NH2	2.17	0.78
27:YD:27:THR:HG21	27:YD:83:GLU:HB3	1.63	0.78
36:YQ:59:ARG:H	36:YQ:59:ARG:HD3	1.48	0.78
1:QA:302:G:O3'	12:QL:17:LYS:HE2	1.83	0.78
11:QK:99:GLN:HG2	11:QK:105:VAL:HG21	1.64	0.78
36:RQ:20:ALA:HB1	36:RQ:99:PRO:HB2	1.65	0.78
3:XC:59:ARG:HH12	3:XC:97:LYS:HE3	1.48	0.78
3:XC:162:GLN:HG2	23:XX:24:A:H1'	1.64	0.78
31:YH:150:ALA:O	31:YH:152:ARG:N	2.14	0.78
33:RN:95:PRO:O	33:RN:97:ARG:N	2.15	0.78
27:YD:94:LEU:HD22	27:YD:95:LEU:N	1.98	0.78
28:YE:137:HIS:HB3	28:YE:138:PRO:HD2	1.65	0.78
31:YH:132:ARG:HB2	31:YH:132:ARG:NH1	1.94	0.78
50:Y4:1:MET:SD	50:Y4:6:HIS:NE2	2.56	0.78
31:RH:153:LYS:CG	31:RH:162:ILE:H	1.96	0.78
33:YN:7:LYS:HD2	33:YN:7:LYS:N	1.94	0.78
5:QE:94:ALA:HB2	5:QE:119:LEU:HG	1.66	0.78
25:YA:2287:A:N6	25:YA:2344:U:H3	1.82	0.78
29:YF:145:GLU:HG3	29:YF:145:GLU:O	1.81	0.78
10:QJ:50:ILE:HA	10:QJ:60:ARG:HG2	1.66	0.78
27:RD:69:ARG:NH2	27:RD:128:GLY:O	2.17	0.78
28:RE:24:THR:HG21	28:RE:188:VAL:CG1	2.13	0.78
50:R4:22:ILE:O	50:R4:24:THR:HG23	1.84	0.78
1:XA:954:G:N2	1:XA:1227:A:H62	1.82	0.78
25:YA:242:G:C5'	54:Y8:62:LEU:CD2	2.61	0.78
27:YD:34:VAL:HG21	27:YD:103:ARG:HA	1.66	0.78
34:YO:88:ASN:HD21	34:YO:92:GLU:HB2	1.47	0.78
25:RA:1542:G:O6	25:RA:1543:A:N6	2.17	0.78
30:RG:6:ALA:N	50:R4:23:GLU:CG	2.46	0.78
27:YD:54:ARG:HG3	27:YD:54:ARG:NH1	1.98	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:4:ILE:HD12	28:YE:28:ALA:HB1	1.67	0.78
25:RA:662:G:OP1	35:RP:15:ARG:NH1	2.17	0.77
10:XJ:57:LYS:HD2	10:XJ:60:ARG:NH2	2.00	0.77
27:YD:25:THR:HG22	27:YD:82:ILE:H	1.46	0.77
31:YH:152:ARG:HG3	31:YH:153:LYS:CE	2.13	0.77
38:YS:106:ARG:HA	38:YS:110:LEU:HD21	1.64	0.77
3:XC:162:GLN:HA	3:XC:162:GLN:HE21	1.46	0.77
10:QJ:50:ILE:CD1	10:QJ:57:LYS:CD	2.63	0.77
28:RE:4:ILE:HD12	28:RE:28:ALA:HB1	1.67	0.77
3:XC:20:SER:HB2	3:XC:40:ARG:HH22	1.50	0.77
13:XM:14:ARG:H	13:XM:44:ARG:HD3	1.49	0.77
27:YD:44:ASN:HD22	27:YD:44:ASN:N	1.79	0.77
25:RA:2107:C:N4	25:RA:2182:G:H1	1.82	0.77
28:RE:55:ASN:C	28:RE:57:LYS:H	1.91	0.77
31:RH:150:ALA:O	31:RH:152:ARG:N	2.14	0.77
50:R4:58:ARG:O	50:R4:63:TYR:HB2	1.84	0.77
25:YA:1009:A:OP2	25:YA:1010:A:OP2	2.02	0.77
1:QA:1298:C:OP2	7:QG:114:ARG:NH2	2.17	0.77
25:RA:1359:A:N1	25:RA:1372:U:C4	2.53	0.77
25:RA:1403:C:H5"	25:RA:1471:A:H1'	1.67	0.77
28:YE:55:ASN:C	28:YE:57:LYS:H	1.91	0.77
29:YF:11:VAL:HB	29:YF:18:ARG:HG3	1.64	0.77
1:QA:1204:A:OP1	14:QN:3:ARG:NH2	2.17	0.77
36:RQ:79:LEU:HD22	36:RQ:79:LEU:C	2.06	0.77
1:XA:1342:C:H4'	9:XI:125:TYR:HB3	1.67	0.77
25:YA:1270:C:H5"	25:YA:1271:G:H5'	1.64	0.77
25:YA:1728:G:N1	25:YA:1730:U:OP2	2.18	0.77
29:YF:20:LEU:HD12	29:YF:21:ALA:H	1.49	0.77
10:QJ:50:ILE:CD1	10:QJ:57:LYS:CG	2.62	0.77
25:RA:265:A:N6	25:RA:427:U:O2'	2.18	0.77
25:YA:2308:G:N1	25:YA:2311:A:H2	1.77	0.77
36:YQ:80:GLU:OE1	46:Y0:7:LEU:HG	1.84	0.77
38:YS:83:LYS:C	38:YS:109:GLY:HA3	2.10	0.77
25:YA:247:G:O6	54:Y8:12:LYS:NZ	2.16	0.77
33:YN:95:PRO:O	33:YN:97:ARG:N	2.18	0.77
54:Y8:59:LYS:CB	54:Y8:59:LYS:HZ3	1.96	0.77
19:QS:5:LEU:HD11	50:R4:67:TYR:CD2	2.19	0.77
19:QS:68:GLY:HA3	50:R4:68:ARG:CB	2.15	0.77
36:RQ:20:ALA:CB	36:RQ:99:PRO:HD2	2.14	0.77
50:R4:1:MET:HB2	50:R4:6:HIS:NE2	2.00	0.77
29:YF:183:VAL:O	29:YF:187:VAL:HG23	1.85	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:YQ:119:ARG:HG2	36:YQ:119:ARG:HH11	1.48	0.77
31:RH:153:LYS:HA	31:RH:153:LYS:NZ	1.99	0.77
25:YA:662:G:OP1	35:YP:15:ARG:NH2	2.18	0.77
27:YD:25:THR:O	27:YD:27:THR:N	2.17	0.77
27:YD:153:ALA:O	27:YD:154:LYS:HG3	1.85	0.77
33:YN:63:THR:CG2	33:YN:66:LYS:NZ	2.47	0.77
5:QE:71:LEU:HD11	5:QE:115:VAL:H	1.49	0.76
31:RH:26:VAL:HG13	31:RH:27:LYS:N	1.96	0.76
36:YQ:20:ALA:CB	36:YQ:99:PRO:HD2	2.14	0.76
36:YQ:20:ALA:HB1	36:YQ:99:PRO:HB2	1.65	0.76
36:YQ:59:ARG:H	36:YQ:59:ARG:CD	1.99	0.76
54:Y8:60:LEU:C	54:Y8:63:PRO:HD2	2.11	0.76
28:RE:14:ILE:HD11	39:RT:14:TYR:CZ	2.19	0.76
30:RG:6:ALA:HB2	50:R4:23:GLU:OE2	1.85	0.76
28:YE:111:ARG:HE	28:YE:160:TYR:HE1	1.31	0.76
29:YF:32:LEU:C	29:YF:32:LEU:HD12	2.10	0.76
36:YQ:79:LEU:HD12	46:Y0:5:LYS:HD2	1.65	0.76
36:YQ:79:LEU:HD22	36:YQ:79:LEU:C	2.07	0.76
36:RQ:66:ILE:HG13	36:RQ:67:ARG:N	1.99	0.76
25:YA:1826:G:H4'	27:YD:242:ARG:NH2	1.99	0.76
38:YS:60:GLY:O	38:YS:61:ASN:HB3	1.83	0.76
1:QA:1002:G:H2'	1:QA:1003:G:H8	1.49	0.76
20:XT:41:ILE:HG22	20:XT:91:LEU:CD1	2.15	0.76
25:YA:1009:A:OP1	33:YN:37:LYS:NZ	2.17	0.76
36:YQ:66:ILE:HG13	36:YQ:67:ARG:N	1.99	0.76
48:Y2:41:ILE:C	48:Y2:41:ILE:HD12	2.10	0.76
1:QA:448:A:OP2	1:QA:485:G:N2	2.18	0.76
36:RQ:90:VAL:HG13	36:RQ:91:GLU:H	1.49	0.76
37:RR:33:ARG:HH22	51:R5:55:ARG:HG2	1.50	0.76
25:YA:574:C:N3	28:YE:145:LYS:NZ	2.34	0.76
31:YH:153:LYS:HA	31:YH:153:LYS:NZ	2.00	0.76
36:YQ:90:VAL:HG13	36:YQ:91:GLU:H	1.49	0.76
37:RR:3:HIS:O	37:RR:5:LYS:N	2.19	0.76
16:XP:45:THR:HG22	16:XP:47:ASP:H	1.50	0.76
25:YA:593:G:H1'	54:Y8:4:MET:HE1	1.68	0.76
25:YA:847:U:C4	25:YA:933:A:N1	2.54	0.76
29:YF:29:ASN:H	29:YF:112:MET:CE	1.97	0.76
25:YA:1247:A:OP2	35:YP:15:ARG:NH1	2.18	0.76
25:YA:2451:A:C6	56:Z6:101:PPU:HE2	2.20	0.76
25:RA:593:G:O2'	54:R8:61:LEU:HD13	1.84	0.76
25:RA:676:A:H8	25:RA:2069:G:H21	1.33	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1359:A:N6	25:RA:1372:U:C4	2.54	0.76
30:RG:5:VAL:HA	50:R4:23:GLU:HG3	1.68	0.76
25:YA:1359:A:C3'	25:YA:1360:A:H5'	2.15	0.76
25:YA:2419:U:O4	54:Y8:30:ARG:CZ	2.33	0.76
25:YA:2849:U:O4	39:YT:23:ARG:NH2	2.18	0.76
25:RA:709:U:H3	25:RA:722:A:H61	1.31	0.76
31:RH:152:ARG:HG3	31:RH:153:LYS:CE	2.13	0.76
31:RH:169:VAL:HG22	31:RH:170:ARG:H	1.48	0.76
32:RI:54:GLN:O	32:RI:58:LEU:HB2	1.85	0.76
51:R5:47:PRO:O	51:R5:48:GLU:HG3	1.86	0.76
31:YH:125:VAL:HA	31:YH:126:PRO:HB3	1.68	0.76
32:YI:144:VAL:HG13	32:YI:145:VAL:HG13	1.67	0.76
38:RS:62:LYS:HB3	38:RS:97:ARG:HD3	1.67	0.76
1:XA:1004:A:H8	1:XA:1036:G:H22	1.34	0.76
25:YA:620:G:H4'	25:YA:621:A:H5"	1.67	0.76
25:YA:780:G:H21	25:YA:783:A:H62	1.34	0.76
28:YE:36:ARG:HH21	28:YE:88:GLY:HA2	1.51	0.76
29:YF:101:LEU:CD1	29:YF:102:PRO:HD2	2.11	0.76
37:YR:74:LYS:O	37:YR:76:VAL:N	2.18	0.76
19:QS:5:LEU:HD13	50:R4:67:TYR:CE2	2.20	0.75
27:YD:69:ARG:HH21	27:YD:130:ALA:CB	1.99	0.75
28:YE:63:LEU:CD1	28:YE:65:GLY:H	1.99	0.75
30:RG:67:LYS:HE2	50:R4:6:HIS:CE1	2.20	0.75
44:RY:95:LYS:HB3	44:RY:100:ALA:HA	1.68	0.75
51:R5:40:LYS:CD	51:R5:46:CYS:HB3	2.15	0.75
54:R8:60:LEU:C	54:R8:63:PRO:HD2	2.11	0.75
1:XA:1321:C:H3'	1:XA:1322:C:H5"	1.67	0.75
4:XD:9:CYS:SG	4:XD:22:LYS:HD2	2.26	0.75
28:YE:23:VAL:HG21	28:YE:183:LEU:HD23	1.69	0.75
54:Y8:59:LYS:HB2	54:Y8:59:LYS:HZ2	1.49	0.75
25:RA:884:C:O2	25:RA:892:G:N1	2.19	0.75
25:RA:1359:A:C6	25:RA:1372:U:C4	2.75	0.75
1:XA:1008:C:H42	1:XA:1021:G:H1	1.31	0.75
20:XT:26:ASN:HB2	20:XT:71:THR:CG2	2.15	0.75
25:RA:2451:A:C6	56:Z5:101:PPU:HE2	2.21	0.75
1:XA:1076:C:H42	1:XA:1081:G:H1	1.34	0.75
4:XD:33:MET:HE2	4:XD:37:PRO:HA	1.68	0.75
25:YA:1093:G:OP1	31:YH:170:ARG:HD2	1.86	0.75
50:R4:34:GLU:HG3	50:R4:35:VAL:H	1.51	0.75
13:XM:93:ARG:NH1	25:YA:888:C:OP1	2.19	0.75
27:YD:142:VAL:HG23	27:YD:193:VAL:HA	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:146:GLU:HB2	27:YD:189:CYS:HB3	1.67	0.75
29:RF:113:ALA:HB1	29:RF:186:ILE:HG21	1.68	0.75
50:R4:71:ARG:HG3	50:R4:71:ARG:NH1	1.90	0.75
31:RH:153:LYS:HG2	31:RH:162:ILE:H	1.52	0.75
44:RY:87:LYS:HD3	44:RY:92:ASN:HB3	1.69	0.75
52:R6:41:PRO:HG2	52:R6:45:LYS:H	1.52	0.75
36:YQ:79:LEU:HD13	46:Y0:5:LYS:HD3	1.68	0.75
39:YT:27:THR:HG23	39:YT:90:GLN:HB3	1.67	0.75
25:RA:49:A:N7	25:RA:120:U:C5	2.55	0.75
25:RA:67:U:H3	25:RA:74:A:H2	1.33	0.75
28:RE:36:ARG:HH21	28:RE:88:GLY:HA2	1.51	0.75
1:QA:686:U:H1'	11:QK:42:TRP:HE1	1.52	0.75
28:RE:63:LEU:CD1	28:RE:65:GLY:H	2.00	0.75
37:RR:56:LYS:NZ	37:RR:90:ARG:O	2.20	0.75
37:RR:74:LYS:O	37:RR:76:VAL:N	2.18	0.75
25:YA:1030:G:OP2	36:YQ:128:LYS:HE2	1.86	0.75
29:YF:7:TYR:HB3	29:YF:21:ALA:CB	2.16	0.75
33:YN:13:TRP:HB2	33:YN:133:GLN:HG3	1.69	0.75
41:YV:24:LYS:HA	41:YV:92:THR:HG23	1.68	0.75
25:RA:2701:C:H3'	25:RA:2702:U:C5'	2.15	0.74
44:RY:86:ARG:HB2	44:RY:95:LYS:HD2	1.69	0.74
47:R1:7:ILE:HG12	47:R1:91:LYS:HZ1	1.52	0.74
28:YE:61:ARG:HB2	28:YE:62:PRO:HD3	1.69	0.74
1:QA:1305:G:H22	1:QA:1331:G:H2'	1.52	0.74
53:R7:9:ARG:HH21	53:R7:48:LYS:CD	1.99	0.74
17:XQ:66:SER:O	17:XQ:70:ARG:NH1	2.21	0.74
25:YA:602:G:O2'	25:YA:604:G:O2'	2.03	0.74
25:YA:1479:G:N7	25:YA:1510:A:N6	2.35	0.74
1:XA:1130:A:O2'	9:XI:3:GLN:NE2	2.20	0.74
27:YD:30:GLU:HG3	27:YD:63:ARG:CZ	2.17	0.74
25:RA:1030:G:OP2	36:RQ:128:LYS:HE2	1.87	0.74
30:YG:161:THR:HG22	30:YG:163:ALA:H	1.53	0.74
36:YQ:79:LEU:O	36:YQ:79:LEU:CD2	2.36	0.74
38:YS:36:TYR:CD2	38:YS:52:SER:HB3	2.23	0.74
39:YT:51:ARG:HG2	39:YT:98:LYS:HG3	1.70	0.74
25:RA:1614:A:H62	42:RW:93:ALA:HB2	1.53	0.74
30:RG:34:LEU:HB2	30:RG:172:LEU:HD21	1.68	0.74
1:XA:1195:C:H5"	1:XA:1196:U:O5'	1.87	0.74
31:YH:150:ALA:C	31:YH:152:ARG:N	2.44	0.74
31:YH:153:LYS:HG2	31:YH:162:ILE:H	1.52	0.74
28:RE:10:GLY:HA3	39:RT:8:LYS:HD2	1.67	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:RI:98:ALA:HB2	32:RI:111:PRO:HB3	1.69	0.74
25:YA:530:G:O2'	25:YA:532:A:N7	2.20	0.74
27:YD:54:ARG:HG3	27:YD:54:ARG:HH11	1.49	0.74
29:YF:129:PHE:HA	29:YF:142:TRP:NE1	2.02	0.74
45:YZ:103:ARG:HB2	45:YZ:138:GLU:HG2	1.69	0.74
28:RE:201:THR:HG22	28:RE:203:LYS:HB3	1.70	0.74
36:RQ:79:LEU:CD1	46:R0:5:LYS:HD3	2.18	0.74
3:XC:162:GLN:HA	3:XC:162:GLN:NE2	2.03	0.74
29:YF:136:THR:HG22	29:YF:166:ALA:O	1.87	0.74
8:QH:29:SER:HB3	8:QH:32:LYS:HG3	1.69	0.74
25:RA:2849:U:OP1	39:RT:95:ARG:NH1	2.20	0.74
30:RG:6:ALA:H	50:R4:23:GLU:HG3	1.53	0.74
50:R4:41:PRO:O	50:R4:42:PHE:HB3	1.87	0.74
25:YA:76:C:O2'	48:Y2:62:THR:HG21	1.88	0.74
25:YA:1403:C:H5"	25:YA:1471:A:H1'	1.70	0.74
28:YE:78:LEU:HG	28:YE:79:ARG:NE	2.03	0.74
1:QA:13:U:O4	1:QA:20:U:O4	2.05	0.74
13:QM:121:LYS:HE2	13:QM:121:LYS:HA	1.68	0.74
25:RA:1607:C:N4	25:RA:1622:G:OP2	2.20	0.74
29:YF:29:ASN:HB3	29:YF:112:MET:HE1	1.70	0.74
5:QE:71:LEU:HD13	5:QE:115:VAL:H	1.52	0.73
25:RA:1061:U:H5'	25:RA:1070:A:H1'	1.68	0.73
36:RQ:79:LEU:O	36:RQ:79:LEU:CD1	2.35	0.73
51:R5:2:ALA:O	51:R5:3:LYS:HB2	1.88	0.73
25:YA:7:G:H1	25:YA:2896:C:H42	1.35	0.73
25:YA:1045:A:C6	25:YA:1111:A:N7	2.56	0.73
31:YH:153:LYS:HG3	31:YH:161:GLY:CA	2.18	0.73
38:YS:62:LYS:HB3	38:YS:97:ARG:HD3	1.68	0.73
43:YX:27:THR:HB	43:YX:80:ILE:HB	1.69	0.73
1:QA:8:A:N6	4:QD:205:GLU:O	2.22	0.73
25:RA:1359:A:N6	25:RA:1372:U:C5	2.55	0.73
28:RE:61:ARG:HB2	28:RE:62:PRO:HD3	1.69	0.73
28:RE:77:ILE:HD12	28:RE:78:LEU:H	1.52	0.73
32:RI:93:THR:HG22	32:RI:119:PRO:HB3	1.69	0.73
50:Y4:9:LEU:H	50:Y4:27:THR:HG23	1.53	0.73
28:RE:203:LYS:HD2	28:RE:203:LYS:O	1.88	0.73
30:RG:3:LEU:CG	50:R4:25:TYR:CE1	2.71	0.73
31:RH:153:LYS:HG3	31:RH:161:GLY:CA	2.18	0.73
39:RT:36:GLU:HG3	39:RT:41:ARG:HD3	1.70	0.73
5:XE:43:LEU:HD22	5:XE:136:MET:CG	2.19	0.73
10:XJ:50:ILE:HD12	10:XJ:57:LYS:HG2	1.68	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:131:LEU:HB2	27:YD:136:ILE:CD1	2.17	0.73
25:RA:1187:G:H5"	41:RV:81:TYR:CE1	2.24	0.73
31:RH:84:SER:O	31:RH:85:LYS:HB2	1.89	0.73
25:YA:2032:G:H21	28:YE:146:THR:HG23	1.52	0.73
27:YD:35:LYS:HZ1	27:YD:65:ILE:HA	1.51	0.73
5:QE:101:ILE:HD13	5:QE:101:ILE:N	2.03	0.73
25:RA:2277:G:H5'	36:RQ:85:LYS:HG3	1.71	0.73
28:RE:23:VAL:HG21	28:RE:183:LEU:HD23	1.68	0.73
1:XA:148:G:H2'	1:XA:149:A:H8	1.53	0.73
54:Y8:61:LEU:O	54:Y8:62:LEU:HB2	1.88	0.73
36:YQ:79:LEU:CD1	36:YQ:79:LEU:O	2.35	0.73
48:Y2:29:LYS:HD3	48:Y2:57:ILE:HD13	1.71	0.73
25:RA:483:A:H4'	44:RY:49:VAL:HA	1.71	0.73
32:RI:4:ILE:HD11	32:RI:44:LEU:HD12	1.69	0.73
4:XD:7:PRO:HB2	4:XD:10:ARG:HD2	1.69	0.73
28:YE:77:ILE:HD12	28:YE:78:LEU:H	1.52	0.73
1:QA:701:C:H1'	1:QA:703:G:C6	2.24	0.73
1:QA:1318:A:H4'	19:QS:11:VAL:HG11	1.69	0.73
3:QC:58:GLU:HB2	3:QC:65:ALA:HB3	1.70	0.73
31:RH:125:VAL:HA	31:RH:126:PRO:HB3	1.68	0.73
29:YF:29:ASN:H	29:YF:112:MET:HE1	1.54	0.73
51:Y5:56:LYS:HG2	51:Y5:58:LEU:HB3	1.70	0.73
3:QC:20:SER:HB2	3:QC:40:ARG:HH22	1.54	0.73
5:QE:69:VAL:CG1	5:QE:71:LEU:HD23	2.19	0.73
25:RA:996:A:H4'	40:RU:92:ARG:HE	1.54	0.73
25:RA:1021:A:H61	25:RA:1142(A):A:H61	1.37	0.73
25:YA:650:C:OP1	54:Y8:48:PHE:CZ	2.42	0.73
25:YA:1309:G:H4'	53:Y7:7:PRO:HB2	1.71	0.73
27:YD:77:ALA:CB	27:YD:97:TYR:HA	2.18	0.73
28:YE:203:LYS:HD2	28:YE:203:LYS:O	1.88	0.73
50:Y4:18:CYS:SG	50:Y4:19:GLY:N	2.62	0.73
28:RE:56:PRO:O	28:RE:57:LYS:HB2	1.89	0.72
36:RQ:90:VAL:CG1	36:RQ:91:GLU:H	2.02	0.72
45:RZ:19:ARG:NH1	45:RZ:84:GLU:O	2.22	0.72
7:QG:9:VAL:HG13	7:QG:94:ARG:HH21	1.54	0.72
31:RH:54:ARG:HH12	31:RH:62:LYS:HG2	1.54	0.72
31:YH:125:VAL:HG12	31:YH:126:PRO:HG3	1.71	0.72
10:QJ:55:LYS:HG3	10:QJ:56:HIS:N	2.05	0.72
25:RA:67:U:N3	25:RA:74:A:C2	2.57	0.72
28:RE:13:ARG:HA	28:RE:22:PRO:HA	1.71	0.72
31:RH:80:SER:O	31:RH:81:GLU:HB2	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:593:G:O2'	54:Y8:61:LEU:CD1	2.35	0.72
36:YQ:90:VAL:CG1	36:YQ:91:GLU:H	2.02	0.72
38:YS:26:LEU:HD23	38:YS:26:LEU:O	1.90	0.72
54:Y8:16:ILE:HD11	54:Y8:57:ARG:HG2	1.69	0.72
1:QA:954:G:H4'	13:QM:121:LYS:HG3	1.71	0.72
51:R5:40:LYS:CE	51:R5:46:CYS:HB3	2.19	0.72
1:XA:1162:C:H42	1:XA:1174:G:H1	1.34	0.72
28:YE:21:VAL:HB	28:YE:22:PRO:CB	2.18	0.72
30:YG:64:THR:HG23	30:YG:66:GLN:H	1.54	0.72
31:YH:54:ARG:HH12	31:YH:62:LYS:HG2	1.54	0.72
31:YH:152:ARG:O	31:YH:153:LYS:HD2	1.90	0.72
54:Y8:29:LYS:HD3	54:Y8:44:LYS:HB2	1.71	0.72
1:QA:692:U:OP1	11:QK:124:LYS:NZ	2.18	0.72
12:QL:126:LYS:HB2	12:QL:126:LYS:NZ	2.04	0.72
28:RE:78:LEU:HG	28:RE:79:ARG:NE	2.03	0.72
31:RH:30:LYS:HD2	31:RH:81:GLU:H	1.54	0.72
1:XA:1108:G:H5'	3:XC:176:HIS:HD1	1.54	0.72
31:YH:132:ARG:HH11	31:YH:132:ARG:CB	1.97	0.72
1:QA:501:C:H2'	1:QA:502:G:H8	1.54	0.72
5:QE:7:GLU:HG2	5:QE:112:LEU:HD22	1.72	0.72
25:RA:2815:C:H5'	51:R5:29:THR:HG21	1.72	0.72
28:RE:28:ALA:HB3	28:RE:93:VAL:HG22	1.72	0.72
12:XL:126:LYS:HB2	12:XL:126:LYS:NZ	2.04	0.72
29:YF:124:LEU:HD12	29:YF:125:LEU:N	2.04	0.72
30:YG:6:ALA:N	50:Y4:23:GLU:HG2	2.04	0.72
35:YP:61:ARG:HD2	54:Y8:13:ARG:CD	2.19	0.72
50:Y4:48:ARG:HH12	50:Y4:52:THR:HG22	1.54	0.72
51:R5:58:LEU:CD1	51:R5:60:VAL:HG12	2.18	0.72
5:XE:91:LEU:HD12	5:XE:120:THR:HG22	1.72	0.72
10:XJ:55:LYS:HG3	10:XJ:56:HIS:N	2.05	0.72
25:YA:242:G:H5'	54:Y8:62:LEU:HD22	1.72	0.72
25:YA:2250:G:C6	36:YQ:82:ARG:HD2	2.25	0.72
28:YE:201:THR:HG22	28:YE:203:LYS:HB3	1.69	0.72
29:YF:32:LEU:HD12	29:YF:32:LEU:O	1.90	0.72
31:RH:26:VAL:CG1	31:RH:27:LYS:H	2.02	0.72
31:RH:89:ILE:CD1	31:RH:129:THR:HB	2.19	0.72
32:RI:3:VAL:HG12	32:RI:38:LEU:HA	1.71	0.72
54:R8:16:ILE:HD11	54:R8:57:ARG:HG2	1.70	0.72
5:XE:10:MET:HB3	5:XE:32:VAL:HG22	1.70	0.72
28:YE:197:ILE:HD11	28:YE:199:ARG:HH12	1.55	0.72
53:Y7:9:ARG:HH21	53:Y7:48:LYS:CD	1.99	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:49:A:N7	25:RA:120:U:C4	2.58	0.72
30:RG:5:VAL:HG22	50:R4:25:TYR:CE2	2.25	0.72
54:R8:61:LEU:O	54:R8:62:LEU:HB2	1.88	0.72
25:YA:2032:G:H21	28:YE:146:THR:CG2	2.03	0.72
28:YE:56:PRO:O	28:YE:57:LYS:HB2	1.89	0.72
29:YF:157:VAL:HB	29:YF:194:MET:HB3	1.70	0.72
38:YS:83:LYS:HZ1	38:YS:109:GLY:HA2	1.50	0.72
44:YY:51:VAL:HG13	44:YY:52:SER:H	1.54	0.72
48:Y2:27:GLU:OE1	48:Y2:27:GLU:N	2.19	0.72
17:QQ:4:LYS:HE3	17:QQ:6:LEU:HD21	1.72	0.72
19:QS:5:LEU:HD22	50:R4:67:TYR:CZ	2.24	0.72
28:RE:197:ILE:HD11	28:RE:199:ARG:HH12	1.55	0.72
31:RH:152:ARG:O	31:RH:153:LYS:HD2	1.90	0.72
36:RQ:79:LEU:O	36:RQ:79:LEU:CD2	2.36	0.72
45:RZ:110:GLY:HA2	45:RZ:111:VAL:C	2.13	0.72
25:YA:617:G:OP1	29:YF:40:GLN:NE2	2.23	0.72
25:YA:957:A:H5'	36:YQ:76:LYS:HD2	1.72	0.72
25:YA:2306:C:H3'	25:YA:2307:G:H5"	1.71	0.72
25:YA:2701:C:H3'	25:YA:2702:U:C5'	2.19	0.72
28:YE:13:ARG:HA	28:YE:22:PRO:HA	1.71	0.72
29:YF:9:ILE:HD11	29:YF:125:LEU:HG	1.70	0.72
31:YH:26:VAL:CG1	31:YH:27:LYS:H	2.02	0.72
25:RA:2135:A:H62	25:RA:2156:G:H21	1.35	0.71
54:R8:60:LEU:O	54:R8:63:PRO:HD2	1.90	0.71
1:XA:272:C:H2'	1:XA:273:A:H8	1.54	0.71
1:XA:1316:G:H22	1:XA:1319:A:H5"	1.55	0.71
1:XA:1450:U:O2'	1:XA:1451:A:N7	2.23	0.71
25:YA:1899:G:H21	25:YA:1902:C:N4	1.88	0.71
25:YA:2636:U:OP1	28:YE:79:ARG:HA	1.90	0.71
34:YO:47:ILE:HG13	34:YO:48:PRO:HD2	1.72	0.71
35:YP:64:LYS:C	35:YP:66:GLY:H	1.97	0.71
38:YS:83:LYS:HG2	38:YS:109:GLY:N	2.04	0.71
38:YS:103:GLU:O	38:YS:106:ARG:HG3	1.90	0.71
2:QB:115:LEU:HB2	2:QB:145:LEU:HD12	1.72	0.71
19:QS:68:GLY:HA3	50:R4:68:ARG:CG	2.20	0.71
28:RE:21:VAL:HB	28:RE:22:PRO:CB	2.18	0.71
28:RE:93:VAL:H	28:RE:95:ILE:HD12	1.54	0.71
46:R0:68:GLU:OE1	46:R0:82:ARG:NH1	2.23	0.71
9:XI:114:TYR:HE2	10:XJ:60:ARG:H	1.38	0.71
31:YH:30:LYS:HD2	31:YH:81:GLU:H	1.54	0.71
31:YH:128:PRO:HD2	31:YH:129:THR:H	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:882:G:H1	25:RA:894:C:H42	0.79	0.71
29:YF:185:ASP:OD1	29:YF:188:ARG:NH1	2.23	0.71
38:YS:67:ARG:O	38:YS:71:ARG:HG3	1.90	0.71
40:YU:92:ARG:HD2	41:YV:11:GLN:HB2	1.72	0.71
54:Y8:60:LEU:O	54:Y8:63:PRO:HD2	1.90	0.71
36:RQ:81:VAL:C	36:RQ:82:ARG:HG2	2.14	0.71
1:XA:148:G:N2	1:XA:174:C:N3	2.34	0.71
1:XA:523:A:H61	12:XL:92:ASP:HB2	1.55	0.71
29:YF:185:ASP:HA	29:YF:188:ARG:CD	2.20	0.71
30:YG:6:ALA:HB2	50:Y4:23:GLU:HG2	1.72	0.71
31:YH:84:SER:O	31:YH:85:LYS:HB2	1.89	0.71
49:R3:8:LEU:HD13	49:R3:31:LEU:HD23	1.71	0.71
1:XA:31:G:O2'	1:XA:48:C:N4	2.22	0.71
38:YS:83:LYS:HZ2	38:YS:109:GLY:HA2	1.55	0.71
54:Y8:58:ILE:HD13	54:Y8:61:LEU:HD11	1.72	0.71
25:RA:1689:A:H62	25:RA:1698:A:H2	1.37	0.71
31:RH:125:VAL:HG12	31:RH:126:PRO:HG3	1.70	0.71
50:R4:29:PRO:O	50:R4:30:GLU:HB2	1.89	0.71
33:YN:89:LYS:O	33:YN:93:THR:HG22	1.90	0.71
36:YQ:81:VAL:C	36:YQ:82:ARG:HG2	2.15	0.71
30:RG:61:ALA:HB2	30:RG:68:PRO:HD3	1.72	0.71
52:R6:17:LYS:HB3	52:R6:44:ARG:HH22	1.55	0.71
8:XH:10:LEU:HD22	8:XH:83:ILE:HD11	1.72	0.71
25:YA:67:U:O4	25:YA:74:A:N1	2.24	0.71
27:YD:263:ARG:NH1	27:YD:263:ARG:HB2	2.06	0.71
31:YH:80:SER:O	31:YH:81:GLU:HB2	1.89	0.71
33:YN:8:GLN:C	33:YN:9:VAL:HG13	2.15	0.71
35:YP:59:LEU:HA	35:YP:61:ARG:HE	1.56	0.71
1:QA:191:G:H1'	20:QT:105:SER:HB3	1.72	0.71
10:QJ:61:GLU:OE2	14:QN:45:ARG:NH1	2.23	0.71
25:RA:1071:G:O6	25:RA:1091:G:O6	2.08	0.71
25:RA:2680:C:H5'	28:RE:189:PRO:HA	1.72	0.71
25:YA:2467:C:H4'	36:YQ:123:HIS:CD2	2.26	0.71
25:YA:2580:U:H4'	28:YE:130:GLY:HA3	1.72	0.71
44:YY:29:GLU:HB3	44:YY:38:ILE:HG23	1.70	0.71
48:Y2:7:ARG:HH11	48:Y2:7:ARG:HG3	1.55	0.71
25:RA:747:U:N1	51:R5:2:ALA:HB3	2.06	0.71
35:RP:95:VAL:HG13	35:RP:100:LEU:HD21	1.73	0.71
54:R8:58:ILE:HD13	54:R8:61:LEU:HD11	1.72	0.71
35:YP:61:ARG:HD2	54:Y8:13:ARG:HD3	1.72	0.71
35:YP:126:VAL:HG13	35:YP:145:PRO:HB2	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:YT:77:PRO:HB2	39:YT:80:SER:HB2	1.72	0.71
44:YY:42:VAL:HG12	44:YY:65:ALA:HB3	1.71	0.71
1:QA:148:G:H1	1:QA:174:C:H42	1.37	0.71
4:QD:32:ALA:O	4:QD:34:GLU:N	2.23	0.71
25:RA:2610:C:H4'	25:RA:2611:U:OP2	1.91	0.71
28:RE:174:ASP:CG	28:RE:175:VAL:H	1.98	0.71
54:R8:59:LYS:CB	54:R8:59:LYS:HZ3	2.02	0.71
1:XA:1002:G:H1	1:XA:1038:C:H42	1.39	0.71
1:XA:1327:C:OP2	21:XU:12:LYS:NZ	2.23	0.71
10:XJ:57:LYS:HD2	10:XJ:60:ARG:CZ	2.21	0.71
12:XL:24:VAL:HG12	12:XL:24:VAL:O	1.90	0.71
25:YA:2445:G:OP1	29:YF:74:ARG:NH2	2.23	0.71
28:YE:93:VAL:H	28:YE:95:ILE:HD12	1.54	0.71
32:YI:5:LEU:HD21	32:YI:12:LEU:HB3	1.72	0.71
34:YO:2:ILE:HD12	34:YO:6:THR:HG21	1.72	0.71
25:RA:1043:C:N4	25:RA:1112:G:H1	1.88	0.70
25:YA:1652:A:OP1	37:YR:8:ARG:NH1	2.24	0.70
28:YE:174:ASP:CG	28:YE:175:VAL:H	1.98	0.70
45:YZ:182:LYS:HG3	45:YZ:183:LEU:HD23	1.70	0.70
19:QS:41:VAL:HB	19:QS:42:PRO:HA	1.73	0.70
25:RA:2392:A:H2	25:RA:2424:C:H42	1.39	0.70
48:R2:29:LYS:HE3	48:R2:57:ILE:HG21	1.73	0.70
29:YF:66:PRO:O	29:YF:67:GLN:HB3	1.89	0.70
31:YH:59:ARG:HH11	31:YH:59:ARG:HG3	1.56	0.70
25:RA:1496:A:H8	25:RA:1577:C:HO2'	1.37	0.70
30:RG:6:ALA:N	50:R4:23:GLU:HG2	2.06	0.70
35:RP:71:VAL:HG13	35:RP:72:PRO:HD3	1.71	0.70
42:RW:29:LEU:HD22	42:RW:69:LEU:HD11	1.72	0.70
27:YD:244:ARG:HB2	27:YD:245:PRO:HD2	1.71	0.70
28:YE:28:ALA:HB3	28:YE:93:VAL:HG22	1.72	0.70
30:YG:47:LYS:HD3	30:YG:81:LYS:HB2	1.72	0.70
31:YH:89:ILE:CD1	31:YH:129:THR:HB	2.20	0.70
31:RH:132:ARG:HH11	31:RH:132:ARG:CB	1.97	0.70
39:RT:18:ASP:OD1	39:RT:18:ASP:N	2.19	0.70
28:YE:14:ILE:HG12	28:YE:15:PHE:N	2.06	0.70
29:YF:178:PRO:HG2	29:YF:179:GLU:OE2	1.90	0.70
31:YH:103:LEU:HD12	31:YH:131:VAL:HG21	1.73	0.70
31:YH:152:ARG:O	31:YH:153:LYS:CB	2.39	0.70
32:YI:3:VAL:HG12	32:YI:38:LEU:HA	1.71	0.70
19:QS:40:ILE:HD11	19:QS:62:ILE:HD12	1.74	0.70
31:RH:154:PRO:O	31:RH:155:SER:HB2	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:RI:54:GLN:O	32:RI:58:LEU:HB3	1.92	0.70
36:RQ:32:TYR:CD1	36:RQ:133:ARG:HA	2.27	0.70
10:XJ:50:ILE:HA	10:XJ:60:ARG:HG2	1.72	0.70
36:YQ:32:TYR:CD1	36:YQ:133:ARG:HA	2.27	0.70
25:RA:1443:G:H1	25:RA:1548:C:N4	1.90	0.70
30:RG:65:GLY:O	50:R4:7:PRO:CD	2.39	0.70
54:R8:29:LYS:HD3	54:R8:44:LYS:HB2	1.71	0.70
10:XJ:61:GLU:OE2	14:XN:45:ARG:NH1	2.25	0.70
25:YA:2680:C:H5'	28:YE:189:PRO:HA	1.74	0.70
29:YF:101:LEU:O	29:YF:106:ARG:NH1	2.23	0.70
29:YF:164:ARG:HG3	29:YF:175:THR:OG1	1.91	0.70
31:YH:86:GLU:CG	31:YH:165:ALA:H	1.94	0.70
25:RA:2308:G:N1	25:RA:2311:A:H2	1.84	0.70
31:RH:154:PRO:HG2	31:RH:162:ILE:O	1.92	0.70
14:XN:48:ALA:HB2	14:XN:53:LEU:HD12	1.73	0.70
27:YD:43:ARG:HB3	27:YD:54:ARG:HB2	1.73	0.70
31:YH:154:PRO:HG2	31:YH:162:ILE:O	1.92	0.70
42:YW:17:VAL:HG12	42:YW:76:VAL:HG11	1.72	0.70
31:RH:150:ALA:C	31:RH:152:ARG:N	2.44	0.70
31:RH:152:ARG:O	31:RH:153:LYS:CB	2.40	0.70
36:RQ:59:ARG:C	36:RQ:60:ARG:HG3	2.16	0.70
18:XR:58:LEU:HD23	18:XR:62:GLU:HB3	1.74	0.70
25:YA:2123:G:H1	25:YA:2175:C:H42	1.39	0.70
28:YE:103:ASP:OD1	28:YE:201:THR:HA	1.92	0.70
29:YF:178:PRO:HB2	29:YF:201:VAL:HG11	1.73	0.70
1:QA:1310:G:OP1	13:QM:77:ASN:ND2	2.25	0.70
25:RA:2682:U:O2'	28:RE:13:ARG:HG2	1.92	0.70
32:RI:144:VAL:C	32:RI:145:VAL:HG12	2.16	0.70
51:R5:40:LYS:HE2	51:R5:47:PRO:HD2	1.73	0.70
2:XB:185:ILE:HG22	2:XB:199:TYR:HB2	1.74	0.70
38:YS:42:ASP:O	38:YS:43:GLU:HB2	1.90	0.70
1:QA:1002:G:H2'	1:QA:1003:G:C8	2.27	0.70
10:QJ:48:THR:HA	10:QJ:62:HIS:HB3	1.72	0.70
13:QM:7:VAL:HG21	30:RG:113:ARG:O	1.92	0.70
31:RH:128:PRO:HD2	31:RH:129:THR:H	1.55	0.70
25:YA:83:G:N2	25:YA:103:A:OP2	2.19	0.70
12:QL:24:VAL:O	12:QL:24:VAL:HG12	1.90	0.69
25:RA:1980:G:O2'	25:RA:1982:C:OP2	2.09	0.69
4:XD:105:VAL:HG13	4:XD:110:PHE:HB2	1.74	0.69
25:YA:2667:C:H1'	31:YH:109:PHE:HD2	1.55	0.69
36:YQ:43:THR:OG1	36:YQ:46:GLN:HB2	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:YS:54:LEU:O	38:YS:54:LEU:HD13	1.91	0.69
13:QM:3:ARG:HH12	30:RG:113:ARG:NH2	1.89	0.69
19:QS:68:GLY:HA3	50:R4:68:ARG:HG2	1.71	0.69
25:RA:1314:C:H42	25:RA:1338:G:H1	1.39	0.69
25:RA:2277:G:P	36:RQ:85:LYS:HB2	2.32	0.69
26:RB:38:C:H42	26:RB:44:G:H1	1.40	0.69
27:RD:93:ALA:HB3	27:RD:105:ILE:HG22	1.74	0.69
1:XA:266:G:C8	1:XA:266:G:H5'	2.28	0.69
1:XA:1023:G:H3'	1:XA:1024:G:H5"	1.72	0.69
20:XT:100:ILE:HG13	20:XT:102:GLY:H	1.56	0.69
25:YA:2584:U:H5'	56:Z6:101:PPU:H103	1.73	0.69
1:QA:1052:U:O2'	1:QA:1055:A:OP2	2.08	0.69
31:RH:59:ARG:HH11	31:RH:59:ARG:HG3	1.56	0.69
19:XS:10:PHE:HB2	19:XS:39:THR:H	1.54	0.69
27:YD:65:ILE:O	27:YD:65:ILE:HD13	1.91	0.69
19:QS:69:HIS:ND1	50:R4:69:LYS:HE2	2.08	0.69
1:XA:1422:G:H5"	34:YO:48:PRO:HB3	1.73	0.69
10:XJ:55:LYS:HG3	10:XJ:56:HIS:CD2	2.27	0.69
25:YA:483:A:H5'	44:YY:49:VAL:HG22	1.74	0.69
48:Y2:47:ASN:O	48:Y2:49:LYS:N	2.25	0.69
1:QA:617:G:H1	1:QA:623:C:H42	1.40	0.69
1:QA:1243:C:OP2	21:QU:10:ARG:NH2	2.25	0.69
28:RE:103:ASP:OD1	28:RE:201:THR:HA	1.92	0.69
31:RH:4:ILE:HG13	31:RH:6:ARG:NE	2.08	0.69
31:RH:103:LEU:HD12	31:RH:131:VAL:HG21	1.73	0.69
45:RZ:94:GLU:HB2	45:RZ:130:PRO:HD2	1.73	0.69
27:YD:65:ILE:HD11	27:YD:67:PHE:CD2	2.27	0.69
29:YF:65:TRP:HZ3	29:YF:73:ALA:O	1.74	0.69
48:Y2:40:SER:C	48:Y2:42:GLY:H	2.00	0.69
28:RE:7:VAL:HG23	28:RE:8:LYS:N	2.07	0.69
28:RE:65:GLY:HA2	28:RE:70:ALA:CB	2.23	0.69
36:RQ:43:THR:OG1	36:RQ:46:GLN:HB2	1.91	0.69
50:R4:37:SER:C	50:R4:39:CYS:H	1.98	0.69
1:XA:1437:C:H42	1:XA:1464:G:H1	1.37	0.69
6:XF:68:PRO:HG2	6:XF:71:ARG:HG3	1.75	0.69
12:XL:25:PRO:C	12:XL:27:LEU:H	1.98	0.69
38:YS:106:ARG:N	38:YS:110:LEU:HD21	2.07	0.69
4:QD:22:LYS:CG	4:QD:26:CYS:SG	2.80	0.69
10:QJ:55:LYS:HG3	10:QJ:56:HIS:CD2	2.27	0.69
22:QV:75:C:H3'	22:QV:76:A:H5"	1.74	0.69
51:R5:4:HIS:HB3	51:R5:5:PRO:HD3	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:538:G:H5"	12:XL:114:LYS:HB2	1.74	0.69
25:YA:1980:G:O2'	25:YA:1982:C:OP2	2.11	0.69
25:YA:2287:A:N6	25:YA:2344:U:N3	2.37	0.69
27:YD:76:PRO:O	27:YD:98:VAL:HG23	1.91	0.69
28:YE:7:VAL:HG23	28:YE:8:LYS:N	2.06	0.69
31:YH:89:ILE:O	31:YH:89:ILE:HG12	1.92	0.69
31:YH:154:PRO:O	31:YH:155:SER:HB2	1.91	0.69
25:RA:807:U:H2'	25:RA:808:G:H8	1.58	0.69
25:RA:2547:U:O2	34:RO:23:ARG:NH2	2.25	0.69
25:RA:2808:U:O4	25:RA:2892:A:N7	2.26	0.69
45:RZ:110:GLY:N	45:RZ:111:VAL:HG12	2.08	0.69
51:R5:40:LYS:HE2	51:R5:47:PRO:CD	2.21	0.69
51:R5:40:LYS:HG2	51:R5:47:PRO:HD2	1.75	0.69
1:XA:1123:A:H4'	10:XJ:36:GLY:HA3	1.75	0.69
27:YD:89:SER:HB2	27:YD:159:ALA:HB2	1.75	0.69
29:YF:67:GLN:O	29:YF:68:LYS:CB	2.39	0.69
52:Y6:11:LEU:HD11	52:Y6:51:GLU:HG3	1.75	0.69
12:QL:25:PRO:C	12:QL:27:LEU:H	1.98	0.69
25:RA:617:G:OP1	29:RF:40:GLN:NE2	2.25	0.69
25:RA:2811:G:P	28:RE:61:ARG:HG3	2.32	0.69
27:RD:25:THR:O	27:RD:27:THR:N	2.26	0.69
32:RI:5:LEU:HD11	32:RI:19:VAL:HG12	1.73	0.69
42:RW:29:LEU:HG	42:RW:33:ARG:HD2	1.74	0.69
25:YA:1826:G:C4'	27:YD:242:ARG:HH21	2.02	0.69
26:YB:44:G:OP1	50:Y4:1:MET:N	2.26	0.69
27:YD:17:THR:HG22	27:YD:205:VAL:N	2.08	0.69
27:YD:122:ASP:CG	27:YD:123:ALA:H	2.00	0.69
40:YU:92:ARG:HG2	40:YU:92:ARG:O	1.91	0.69
40:YU:92:ARG:NH1	41:YV:11:GLN:O	2.26	0.69
25:RA:2114:A:N6	25:RA:2119:A:N7	2.41	0.69
36:RQ:80:GLU:HG3	36:RQ:81:VAL:H	1.58	0.69
27:YD:35:LYS:HB3	27:YD:63:ARG:HA	1.75	0.69
1:QA:1312:G:H5"	50:R4:67:TYR:OH	1.93	0.68
25:RA:2712:U:HO2'	25:RA:2712(A):A:H8	1.38	0.68
32:RI:41:GLU:HA	32:RI:44:LEU:HB2	1.74	0.68
45:RZ:60:GLU:HA	45:RZ:66:SER:HA	1.75	0.68
25:YA:660:G:O3'	29:YF:38:ARG:NH2	2.26	0.68
27:YD:17:THR:CG2	27:YD:204:ILE:HA	2.23	0.68
35:YP:62:LEU:CD2	54:Y8:25:MET:HB2	2.23	0.68
45:YZ:60:GLU:HA	45:YZ:66:SER:HA	1.75	0.68
47:Y1:73:LEU:HD13	47:Y1:90:ILE:HG22	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:49:U:C4	1:QA:365:U:O4	2.45	0.68
25:RA:242:G:H5'	54:R8:62:LEU:HD22	1.73	0.68
25:RA:2392:A:C8	35:RP:60:MET:HG2	2.28	0.68
1:XA:503:C:OP2	12:XL:116:SER:HB3	1.94	0.68
2:XB:74:LYS:HE3	2:XB:166:ASP:CB	2.22	0.68
38:YS:57:LYS:HD3	38:YS:57:LYS:H	1.58	0.68
1:XA:1077:G:N2	1:XA:1080:A:OP2	2.23	0.68
28:YE:65:GLY:HA2	28:YE:70:ALA:CB	2.23	0.68
31:YH:126:PRO:HB2	31:YH:130:ARG:O	1.93	0.68
35:YP:19:VAL:HG13	35:YP:21:ARG:H	1.57	0.68
38:YS:106:ARG:CA	38:YS:110:LEU:HD11	2.19	0.68
1:QA:589:C:N4	1:QA:650:G:H1	1.89	0.68
3:QC:3:ASN:N	3:QC:3:ASN:OD1	2.26	0.68
25:RA:2821:A:OP2	28:RE:110:GLY:HA3	1.93	0.68
25:YA:2753:A:O2'	55:Y9:15:LYS:NZ	2.27	0.68
31:YH:4:ILE:HG13	31:YH:6:ARG:NE	2.08	0.68
36:YQ:88:GLY:C	36:YQ:90:VAL:N	2.47	0.68
38:YS:100:ALA:HA	38:YS:103:GLU:HG2	1.75	0.68
39:YT:123:GLN:O	39:YT:125:ARG:N	2.26	0.68
30:RG:47:LYS:HD3	30:RG:81:LYS:HB2	1.76	0.68
30:RG:66:GLN:NE2	30:RG:93:THR:O	2.26	0.68
25:YA:141:A:H8	25:YA:1408:C:H1'	1.58	0.68
25:YA:2600:A:N7	27:YD:237:GLU:OE2	2.27	0.68
36:YQ:80:GLU:HG3	36:YQ:81:VAL:H	1.58	0.68
37:YR:33:ARG:NH2	51:Y5:55:ARG:HG2	2.08	0.68
1:QA:542:G:OP1	4:QD:10:ARG:NH2	2.26	0.68
31:RH:126:PRO:HB2	31:RH:130:ARG:O	1.93	0.68
14:XN:13:THR:N	14:XN:14:PRO:HD2	2.09	0.68
28:YE:15:PHE:CE1	39:YT:81:PRO:HD2	2.28	0.68
31:YH:4:ILE:HG13	31:YH:6:ARG:NH1	2.09	0.68
31:YH:126:PRO:HG2	31:YH:127:GLU:H	1.58	0.68
25:RA:530:G:C2	25:RA:2022:U:OP1	2.46	0.68
36:RQ:12:GLN:CG	36:RQ:73:PRO:HD2	2.21	0.68
36:RQ:88:GLY:C	36:RQ:90:VAL:N	2.47	0.68
50:R4:15:ILE:HD13	50:R4:15:ILE:N	2.09	0.68
2:XB:235:SER:OG	2:XB:236:TYR:N	2.25	0.68
25:YA:2278:A:OP1	36:YQ:11:LYS:HD2	1.93	0.68
36:YQ:66:ILE:HG13	36:YQ:67:ARG:H	1.58	0.68
5:QE:11:ILE:HG13	5:QE:31:LEU:HB3	1.76	0.68
16:QP:4:ILE:HG12	16:QP:21:VAL:HG12	1.75	0.68
31:RH:88:LEU:H	31:RH:88:LEU:HD22	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:RH:126:PRO:HG2	31:RH:127:GLU:H	1.59	0.68
25:YA:1845:G:OP1	27:YD:258:LYS:NZ	2.23	0.68
27:YD:241:PRO:O	27:YD:243:GLY:N	2.27	0.68
28:YE:16:ARG:HG3	28:YE:16:ARG:O	1.93	0.68
31:YH:126:PRO:CD	31:YH:127:GLU:H	2.07	0.68
35:YP:88:LEU:HD12	35:YP:95:VAL:HG11	1.76	0.68
45:YZ:145:GLU:HG3	45:YZ:146:ILE:HG12	1.76	0.68
48:Y2:23:LYS:O	48:Y2:27:GLU:OE1	2.11	0.68
37:RR:38:VAL:HG22	37:RR:112:ALA:HB2	1.75	0.68
51:R5:20:ARG:HA	51:R5:23:HIS:ND1	2.09	0.68
19:XS:40:ILE:HG12	19:XS:41:VAL:HG13	1.75	0.68
28:YE:9:VAL:HB	28:YE:25:VAL:HG23	1.76	0.68
29:YF:34:TRP:CE3	35:YP:8:PRO:HB3	2.29	0.68
33:YN:133:GLN:HB2	33:YN:135:PRO:HD3	1.76	0.68
35:YP:65:ARG:HH21	54:Y8:46:ARG:HH12	1.40	0.68
5:QE:71:LEU:CD1	5:QE:115:VAL:N	2.54	0.68
11:QK:58:PRO:HB2	11:QK:93:GLN:HG3	1.75	0.68
25:RA:1693:U:O2'	27:RD:14:ARG:NH2	2.27	0.68
30:RG:83:ARG:H	30:RG:86:MET:HG3	1.58	0.68
36:RQ:104:PHE:HE2	36:RQ:125:LEU:HD11	1.59	0.68
25:YA:2789:C:H1'	25:YA:2892:A:H2	1.59	0.68
38:YS:35:ILE:HD13	38:YS:101:LEU:HD23	1.76	0.68
1:QA:1175:G:H2'	1:QA:1176:A:H8	1.59	0.67
1:QA:1305:G:N2	1:QA:1331:G:H2'	2.09	0.67
5:QE:71:LEU:CD2	5:QE:115:VAL:HG22	2.24	0.67
25:RA:2584:U:H5'	56:Z5:101:PPU:H103	1.77	0.67
25:RA:2683:C:OP1	39:RT:53:ARG:NH2	2.27	0.67
28:RE:116:VAL:O	28:RE:117:MET:HB3	1.94	0.67
27:YD:44:ASN:N	27:YD:44:ASN:ND2	2.42	0.67
36:YQ:90:VAL:CG1	36:YQ:91:GLU:N	2.57	0.67
48:Y2:64:LEU:HD22	48:Y2:68:ARG:HD2	1.77	0.67
1:QA:1502:A:H2	1:QA:1505:G:H1	1.41	0.67
25:RA:2580:U:H4'	28:RE:130:GLY:HA3	1.77	0.67
28:RE:13:ARG:CB	28:RE:13:ARG:HH11	2.07	0.67
36:RQ:20:ALA:HB1	36:RQ:99:PRO:HD2	1.76	0.67
36:RQ:133:ARG:O	36:RQ:134:ARG:HB2	1.94	0.67
24:XY:30:C:H42	24:XY:40:G:H1	1.41	0.67
31:YH:77:LYS:HG2	31:YH:77:LYS:O	1.94	0.67
36:YQ:133:ARG:O	36:YQ:134:ARG:HB2	1.94	0.67
1:QA:376:G:H1	1:QA:387:U:H3	1.42	0.67
1:QA:1347:G:N2	1:QA:1374:A:OP2	2.20	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:301:G:H1	25:RA:316:C:N4	1.89	0.67
30:RG:6:ALA:HB2	50:R4:23:GLU:HG2	1.73	0.67
31:RH:89:ILE:O	31:RH:89:ILE:HG12	1.93	0.67
36:RQ:90:VAL:CG1	36:RQ:91:GLU:N	2.57	0.67
36:YQ:12:GLN:CG	36:YQ:73:PRO:HD2	2.21	0.67
36:YQ:90:VAL:O	36:YQ:92:GLY:N	2.25	0.67
45:YZ:19:ARG:NH1	45:YZ:84:GLU:O	2.24	0.67
1:QA:1286:A:H5"	21:QU:26:LYS:HD2	1.75	0.67
25:RA:414:C:O2	25:RA:1864:U:O2'	2.13	0.67
26:RB:48:A:H4'	38:RS:95:HIS:HD2	1.60	0.67
28:RE:10:GLY:H	28:RE:25:VAL:HG23	1.59	0.67
25:YA:1812:A:O2'	27:YD:45:ASN:HB2	1.95	0.67
36:YQ:33:GLY:HA2	36:YQ:105:GLU:HA	1.76	0.67
38:YS:26:LEU:HD12	38:YS:39:ILE:CD1	2.23	0.67
38:YS:52:SER:O	38:YS:56:LEU:HD22	1.93	0.67
38:YS:67:ARG:HB2	38:YS:67:ARG:CZ	2.24	0.67
51:Y5:48:GLU:HA	51:Y5:59:GLU:CG	2.24	0.67
2:QB:27:LYS:HD2	2:QB:193:ASP:HB2	1.75	0.67
25:RA:259:G:H21	25:RA:621:A:H8	1.40	0.67
28:RE:14:ILE:HG12	28:RE:15:PHE:N	2.06	0.67
50:R4:33:VAL:HG12	50:R4:34:GLU:N	2.10	0.67
53:R7:9:ARG:NH2	53:R7:48:LYS:HB2	2.10	0.67
1:XA:1502:A:H2	1:XA:1505:G:H1	1.42	0.67
8:XH:120:THR:H	8:XH:123:GLU:HB2	1.59	0.67
28:YE:116:VAL:O	28:YE:117:MET:HB3	1.94	0.67
36:YQ:104:PHE:CE2	36:YQ:125:LEU:HD11	2.29	0.67
53:Y7:9:ARG:CZ	53:Y7:48:LYS:HB2	2.23	0.67
1:QA:1162:C:H42	1:QA:1174:G:H1	1.42	0.67
19:QS:29:ARG:HD3	19:QS:30:LEU:HD13	1.77	0.67
36:RQ:104:PHE:CE2	36:RQ:125:LEU:HD11	2.30	0.67
41:RV:72:VAL:HG13	41:RV:85:LYS:HB3	1.75	0.67
5:XE:31:LEU:HD23	5:XE:45:PHE:CD1	2.30	0.67
28:YE:26:ILE:HD13	28:YE:27:LEU:N	2.10	0.67
25:RA:270(T):G:H5"	47:R1:97:LEU:HD22	1.75	0.67
25:RA:1226:G:H4'	41:RV:84:LYS:HG2	1.77	0.67
25:RA:2839:G:H5'	37:RR:46:GLY:HA2	1.76	0.67
27:RD:182:LEU:H	27:RD:272:ALA:HB3	1.59	0.67
31:RH:4:ILE:HG13	31:RH:6:ARG:NH1	2.09	0.67
39:RT:84:GLN:HG2	39:RT:85:LYS:HG2	1.76	0.67
25:YA:1607:C:N4	25:YA:1622:G:OP2	2.27	0.67
28:YE:13:ARG:CB	28:YE:13:ARG:HH11	2.07	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:YH:88:LEU:H	31:YH:88:LEU:HD22	1.58	0.67
41:YV:52:VAL:HG21	41:YV:55:ALA:HB3	1.76	0.67
19:QS:5:LEU:CD2	50:R4:67:TYR:OH	2.42	0.67
25:RA:857:C:H4'	46:R0:23:VAL:HG21	1.75	0.67
30:RG:3:LEU:HD21	50:R4:25:TYR:CE1	2.29	0.67
36:RQ:32:TYR:HD1	36:RQ:133:ARG:HA	1.60	0.67
36:RQ:90:VAL:O	36:RQ:92:GLY:N	2.25	0.67
37:RR:33:ARG:NH2	51:R5:55:ARG:CG	2.58	0.67
2:XB:168:THR:HB	2:XB:192:SER:HB2	1.77	0.67
25:YA:1258:C:O4'	29:YF:84:VAL:HG11	1.95	0.67
29:YF:107:LYS:O	29:YF:108:LYS:C	2.37	0.67
29:YF:184:TYR:O	29:YF:188:ARG:HG3	1.94	0.67
25:RA:747:U:C2	51:R5:2:ALA:HB3	2.29	0.67
25:RA:958:U:OP2	36:RQ:14:ARG:NH1	2.27	0.67
25:RA:1309:G:H4'	53:R7:7:PRO:HB2	1.75	0.67
28:RE:16:ARG:HG3	28:RE:16:ARG:O	1.93	0.67
31:RH:124:GLU:HB3	31:RH:132:ARG:HD2	1.77	0.67
31:RH:125:VAL:CG1	31:RH:126:PRO:HG3	2.25	0.67
31:RH:126:PRO:CD	31:RH:127:GLU:H	2.07	0.67
36:RQ:66:ILE:HG13	36:RQ:67:ARG:H	1.57	0.67
2:XB:174:VAL:HG13	2:XB:184:VAL:HG11	1.76	0.67
25:YA:1794:U:H2'	25:YA:1795:C:H6	1.58	0.67
25:YA:2123:G:H2'	25:YA:2124:G:H8	1.60	0.67
28:YE:62:PRO:O	28:YE:64:LYS:N	2.28	0.67
39:YT:16:ARG:NH2	39:YT:83:ILE:O	2.27	0.67
1:QA:1124:G:H3'	1:QA:1145:C:N4	2.10	0.67
27:RD:49:ILE:HD11	27:RD:52:ARG:HA	1.77	0.67
5:XE:100:VAL:HG22	5:XE:118:ILE:HG22	1.77	0.67
36:YQ:104:PHE:HE2	36:YQ:125:LEU:HD11	1.59	0.67
37:YR:78:LYS:HE2	37:YR:83:ILE:HD11	1.77	0.67
1:QA:1055:A:O2'	3:QC:161:GLU:OE2	2.14	0.66
25:RA:1454:U:OP1	37:RR:77:ARG:NH1	2.27	0.66
31:RH:77:LYS:O	31:RH:77:LYS:HG2	1.94	0.66
1:XA:1346:A:OP1	9:XI:120:ARG:NH1	2.26	0.66
27:YD:35:LYS:CA	27:YD:64:ILE:HG22	2.26	0.66
28:YE:51:PHE:O	28:YE:52:LEU:C	2.38	0.66
29:YF:46:ARG:HH11	29:YF:46:ARG:CG	2.04	0.66
29:YF:103:LYS:HA	29:YF:106:ARG:CG	2.21	0.66
25:RA:593:G:H1	25:RA:664:C:H42	1.42	0.66
28:RE:36:ARG:HB3	28:RE:36:ARG:HH11	1.60	0.66
30:RG:6:ALA:N	50:R4:23:GLU:HG3	2.08	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:1095:U:P	1:XA:1108:G:H1	2.18	0.66
25:YA:712:G:H1	25:YA:719:C:H42	1.43	0.66
25:YA:1348:G:H2'	25:YA:1349:A:H5"	1.77	0.66
25:YA:2277:G:C5'	36:YQ:85:LYS:HG3	2.25	0.66
27:YD:35:LYS:HZ1	27:YD:104:TYR:HB2	1.58	0.66
32:YI:4:ILE:HG12	32:YI:18:VAL:HG22	1.77	0.66
32:YI:8:PRO:HD3	32:YI:15:VAL:HG13	1.77	0.66
36:YQ:32:TYR:HD1	36:YQ:133:ARG:HA	1.60	0.66
5:QE:107:ARG:HG3	5:QE:108:ALA:N	2.11	0.66
25:YA:888:C:H3'	25:YA:889:C:H4'	1.77	0.66
25:YA:1332:G:N2	25:YA:1609:A:O2'	2.28	0.66
48:Y2:65:ASN:HB3	48:Y2:69:ARG:NH1	2.10	0.66
1:QA:589:C:N3	1:QA:650:G:N2	2.35	0.66
25:RA:1064:C:N3	25:RA:1074:G:N2	2.40	0.66
25:RA:2620:C:OP1	28:RE:152:LYS:O	2.14	0.66
55:R9:27:CYS:SG	55:R9:29:ASN:ND2	2.69	0.66
1:XA:790:A:OP1	22:XV:38:A:O2'	2.10	0.66
28:YE:28:ALA:O	28:YE:93:VAL:HG23	1.96	0.66
38:YS:106:ARG:HA	38:YS:110:LEU:CD2	2.25	0.66
42:YW:45:TYR:CZ	42:YW:49:LYS:HD2	2.30	0.66
3:QC:162:GLN:HA	3:QC:162:GLN:HE21	1.59	0.66
5:QE:76:ILE:HG23	5:QE:118:ILE:HD13	1.78	0.66
25:RA:1264:G:H5'	51:R5:11:THR:HG21	1.75	0.66
25:RA:2123:G:H2'	25:RA:2124:G:H8	1.60	0.66
31:RH:168:PRO:O	31:RH:169:VAL:HG12	1.96	0.66
36:RQ:86:GLY:C	36:RQ:88:GLY:H	2.03	0.66
13:XM:105:THR:O	13:XM:107:ALA:N	2.29	0.66
25:YA:2811:G:P	28:YE:61:ARG:HG3	2.34	0.66
27:YD:68:LYS:HB2	27:YD:70:TRP:CZ3	2.31	0.66
27:YD:135:PHE:N	27:YD:135:PHE:CD1	2.62	0.66
27:YD:172:TYR:HB3	27:YD:184:LYS:HG2	1.77	0.66
28:YE:37:ARG:NE	28:YE:37:ARG:HA	2.11	0.66
36:YQ:79:LEU:CD1	46:Y0:5:LYS:CD	2.62	0.66
3:QC:70:VAL:HG12	3:QC:72:LYS:H	1.60	0.66
5:QE:101:ILE:HD11	5:QE:119:LEU:HD23	1.78	0.66
25:RA:270(G):C:N4	25:RA:270(S):G:H1	1.94	0.66
25:RA:586:A:H5'	29:RF:89:VAL:HG21	1.77	0.66
25:RA:993:G:OP1	40:RU:50:ARG:NH2	2.28	0.66
25:RA:2635:C:OP1	28:RE:78:LEU:HD12	1.95	0.66
28:RE:9:VAL:HB	28:RE:25:VAL:HG23	1.76	0.66
32:RI:133:HIS:HB2	32:RI:134:PRO:HD2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:R1:7:ILE:HG12	47:R1:91:LYS:NZ	2.11	0.66
1:XA:690:G:H22	11:XK:55:LYS:NZ	1.92	0.66
10:XJ:7:LYS:HB2	10:XJ:97:GLU:HB2	1.76	0.66
25:YA:2391:G:OP2	54:Y8:32:LEU:HD13	1.96	0.66
44:YY:49:VAL:O	44:YY:51:VAL:N	2.29	0.66
19:QS:39:THR:HG22	19:QS:40:ILE:H	1.61	0.66
25:RA:884:C:N3	25:RA:892:G:O6	2.29	0.66
25:RA:2112:G:O6	25:RA:2169:A:N6	2.29	0.66
28:RE:101:ARG:CZ	28:RE:171:GLU:HB2	2.25	0.66
36:RQ:33:GLY:HA2	36:RQ:105:GLU:HA	1.76	0.66
12:XL:26:ALA:O	12:XL:27:LEU:O	2.14	0.66
1:QA:411:A:H62	1:QA:413:G:H21	1.44	0.66
19:QS:5:LEU:HD11	50:R4:67:TYR:CE2	2.30	0.66
25:RA:1332:G:H21	25:RA:1610:A:H8	1.41	0.66
28:RE:62:PRO:O	28:RE:64:LYS:N	2.28	0.66
27:YD:35:LYS:HE3	27:YD:64:ILE:C	2.21	0.66
31:YH:125:VAL:CG1	31:YH:126:PRO:HG3	2.25	0.66
44:YY:97:ARG:HE	44:YY:98:VAL:HB	1.61	0.66
54:Y8:30:ARG:O	54:Y8:31:HIS:HB2	1.96	0.66
25:RA:2233:U:H2'	25:RA:2234:G:C8	2.30	0.66
44:RY:49:VAL:O	44:RY:51:VAL:N	2.29	0.66
25:YA:2304:G:H22	25:YA:2312:U:H3	1.42	0.66
31:YH:124:GLU:HB3	31:YH:132:ARG:HD2	1.77	0.66
36:YQ:20:ALA:HB1	36:YQ:99:PRO:HD2	1.77	0.66
36:YQ:88:GLY:C	36:YQ:90:VAL:H	2.02	0.66
38:YS:88:ASP:OD1	38:YS:90:GLY:N	2.28	0.66
12:QL:126:LYS:C	12:QL:128:ALA:H	2.03	0.66
25:RA:2651:C:H42	25:RA:2669:G:H1	1.43	0.66
28:RE:26:ILE:HD13	28:RE:27:LEU:N	2.10	0.66
50:R4:16:CYS:SG	50:R4:33:VAL:HB	2.35	0.66
19:XS:50:ALA:HB1	19:XS:57:HIS:HB3	1.77	0.66
25:YA:2277:G:H5'	36:YQ:85:LYS:HG3	1.76	0.66
33:YN:35:ARG:O	33:YN:37:LYS:N	2.29	0.66
38:YS:107:GLU:H	38:YS:110:LEU:HD11	1.60	0.66
41:YV:21:ARG:HD2	41:YV:91:TYR:CD1	2.31	0.66
20:QT:26:ASN:HB2	20:QT:71:THR:HG23	1.78	0.65
39:RT:102:ILE:HB	39:RT:110:ILE:HD13	1.78	0.65
1:XA:1124:G:H3'	1:XA:1145:C:N4	2.11	0.65
1:XA:1368:G:OP1	9:XI:111:ARG:NH2	2.25	0.65
22:XV:5:G:H1	22:XV:67:C:N4	1.94	0.65
27:YD:121:PRO:HB3	27:YD:135:PHE:CE2	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:13:ARG:NH1	28:YE:21:VAL:HG12	2.11	0.65
33:YN:1:MET:HE1	40:YU:95:LEU:HD21	1.76	0.65
47:Y1:29:GLY:O	47:Y1:31:GLY:N	2.30	0.65
25:RA:2015:A:N3	51:R5:2:ALA:N	2.44	0.65
25:YA:1689:A:H62	25:YA:1698:A:H2	1.43	0.65
27:YD:183:ARG:HH11	27:YD:183:ARG:CG	2.07	0.65
39:YT:16:ARG:HD3	39:YT:19:LEU:HD11	1.77	0.65
12:QL:115:LYS:O	12:QL:117:ARG:HG3	1.96	0.65
27:RD:108:PRO:HG2	27:RD:111:LEU:HG	1.79	0.65
12:XL:39:VAL:HB	12:XL:57:LYS:HB2	1.78	0.65
25:YA:1533:C:H42	25:YA:1538:G:H1	1.44	0.65
28:YE:36:ARG:HB3	28:YE:36:ARG:HH11	1.60	0.65
28:YE:101:ARG:CZ	28:YE:171:GLU:HB2	2.26	0.65
33:YN:8:GLN:O	33:YN:9:VAL:HG22	1.96	0.65
45:YZ:97:GLU:HB3	45:YZ:125:LEU:HD11	1.78	0.65
1:QA:791:G:H2'	1:QA:792:A:H5'	1.79	0.65
2:QB:82:ARG:HA	2:QB:92:TYR:CE2	2.31	0.65
12:QL:21:LYS:HD2	12:QL:21:LYS:N	2.11	0.65
25:RA:2010:G:H5"	42:RW:42:ARG:HB2	1.79	0.65
27:RD:27:THR:HG21	27:RD:81:ALA:HB1	1.78	0.65
28:RE:11:MET:HE3	28:RE:186:GLY:HA2	1.79	0.65
1:XA:973:G:H3'	1:XA:974:A:H5"	1.77	0.65
22:XV:14:G:H22	22:XV:48:C:H42	1.42	0.65
25:YA:2023:G:H5'	25:YA:2617:C:H4'	1.78	0.65
25:YA:2511:U:O4	25:YA:2575:C:N3	2.30	0.65
27:YD:80:ALA:HB3	27:YD:94:LEU:CD1	2.26	0.65
13:QM:50:GLU:OE1	50:R4:32:TYR:CE2	2.50	0.65
25:RA:593:G:O3'	54:R8:61:LEU:HD22	1.97	0.65
25:RA:2306:C:H3'	25:RA:2307:G:H5"	1.77	0.65
28:RE:28:ALA:O	28:RE:93:VAL:HG23	1.95	0.65
36:RQ:88:GLY:C	36:RQ:90:VAL:H	2.02	0.65
51:R5:40:LYS:NZ	51:R5:48:GLU:HB2	2.10	0.65
52:R6:11:LEU:HD23	52:R6:26:ASN:HB3	1.78	0.65
20:XT:49:ALA:HB1	20:XT:99:LEU:HB2	1.77	0.65
25:YA:277:C:H5'	25:YA:278:A:H5"	1.77	0.65
25:YA:1496:A:H8	25:YA:1577:C:HO2'	1.44	0.65
27:YD:27:THR:CG2	27:YD:28:GLU:H	2.08	0.65
28:YE:201:THR:HG22	28:YE:203:LYS:N	2.07	0.65
40:YU:8:VAL:HG23	40:YU:11:ARG:HH21	1.62	0.65
48:Y2:42:GLY:O	48:Y2:44:LEU:N	2.30	0.65
48:Y2:69:ARG:NH1	48:Y2:69:ARG:HB2	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:Y5:40:LYS:HZ1	51:Y5:48:GLU:HB2	1.61	0.65
19:QS:28:LYS:HB2	19:QS:47:HIS:CE1	2.32	0.65
25:RA:747:U:C6	51:R5:2:ALA:O	2.49	0.65
36:RQ:86:GLY:C	36:RQ:88:GLY:N	2.52	0.65
50:R4:49:PHE:O	50:R4:50:VAL:HG23	1.97	0.65
20:XT:97:ALA:O	20:XT:99:LEU:N	2.30	0.65
25:YA:1754:C:OP1	39:YT:96:ARG:NH1	2.30	0.65
25:YA:2646:C:OP2	25:YA:2732:G:O2'	2.15	0.65
27:YD:145:VAL:HG12	27:YD:146:GLU:O	1.96	0.65
31:YH:168:PRO:O	31:YH:169:VAL:HG12	1.96	0.65
3:QC:9:GLY:HA2	3:QC:12:LEU:HD23	1.78	0.65
16:QP:53:VAL:HG12	16:QP:79:VAL:HG22	1.77	0.65
19:QS:68:GLY:HA2	50:R4:68:ARG:HG2	1.79	0.65
25:RA:141:A:H8	25:RA:1595:G:H21	1.44	0.65
28:RE:51:PHE:O	28:RE:52:LEU:C	2.38	0.65
31:RH:128:PRO:CD	31:RH:129:THR:H	2.09	0.65
35:RP:19:VAL:HG13	35:RP:21:ARG:H	1.61	0.65
36:RQ:80:GLU:C	36:RQ:81:VAL:HG13	2.22	0.65
44:RY:38:ILE:HG22	44:RY:66:PRO:HA	1.79	0.65
50:R4:36:CYS:O	50:R4:37:SER:O	2.14	0.65
2:XB:178:ARG:NH1	2:XB:196:LEU:O	2.29	0.65
10:XJ:50:ILE:HD11	10:XJ:57:LYS:CG	2.25	0.65
12:XL:115:LYS:O	12:XL:117:ARG:HG3	1.97	0.65
22:XV:75:C:OP1	25:YA:2602:A:OP1	2.15	0.65
28:YE:10:GLY:H	28:YE:25:VAL:HG23	1.60	0.65
45:YZ:58:VAL:O	45:YZ:60:GLU:N	2.29	0.65
1:QA:671:G:N2	1:QA:735:C:N3	2.43	0.65
12:QL:39:VAL:HB	12:QL:57:LYS:HB2	1.79	0.65
28:RE:4:ILE:C	28:RE:5:LEU:HD23	2.22	0.65
46:Y0:10:THR:HG22	46:Y0:12:ASN:H	1.62	0.65
51:Y5:48:GLU:HA	51:Y5:59:GLU:HG3	1.78	0.65
51:Y5:56:LYS:H	51:Y5:56:LYS:HD2	1.60	0.65
13:QM:3:ARG:HD2	13:QM:9:ILE:HG12	1.79	0.65
25:RA:2250:G:C6	36:RQ:82:ARG:HD2	2.32	0.65
28:RE:37:ARG:NE	28:RE:37:ARG:HA	2.11	0.65
1:XA:1451:A:H5"	1:XA:1452:C:H5'	1.78	0.65
11:XK:21:ILE:HB	11:XK:84:VAL:HG12	1.78	0.65
25:YA:482:A:H4'	44:YY:47:LYS:HD2	1.78	0.65
27:YD:77:ALA:HB2	27:YD:97:TYR:HA	1.77	0.65
27:YD:176:ARG:HG2	27:YD:176:ARG:HH11	1.61	0.65
54:Y8:52:LYS:HG3	54:Y8:52:LYS:O	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:QM:3:ARG:CZ	30:RG:113:ARG:HH21	2.10	0.65
19:QS:5:LEU:CD2	50:R4:67:TYR:CE2	2.79	0.65
25:RA:631:A:OP2	54:R8:46:ARG:NH2	2.28	0.65
31:RH:51:ARG:HH11	31:RH:51:ARG:HG3	1.61	0.65
51:R5:40:LYS:HD3	51:R5:46:CYS:CB	2.26	0.65
1:XA:1178:G:N2	1:XA:1181:G:N7	2.45	0.65
5:XE:102:ALA:HB2	5:XE:120:THR:OG1	1.97	0.65
12:XL:21:LYS:HD2	12:XL:21:LYS:N	2.11	0.65
26:YB:44:G:H1'	26:YB:47:C:H42	1.62	0.65
30:YG:6:ALA:CB	50:Y4:23:GLU:HG2	2.27	0.65
31:YH:128:PRO:CD	31:YH:129:THR:H	2.09	0.65
54:Y8:56:GLU:OE1	54:Y8:56:GLU:N	2.30	0.65
12:QL:26:ALA:O	12:QL:27:LEU:O	2.14	0.64
25:RA:1224:G:N2	25:RA:1227:A:OP2	2.26	0.64
25:RA:1754:C:OP1	39:RT:96:ARG:NH1	2.25	0.64
26:RB:75:G:H5"	45:RZ:36:LYS:HE2	1.79	0.64
28:RE:104:VAL:HG11	28:RE:188:VAL:CG2	2.27	0.64
31:RH:105:LEU:H	31:RH:105:LEU:CD1	2.09	0.64
51:R5:56:LYS:H	51:R5:56:LYS:CD	2.07	0.64
1:XA:110:C:O2'	16:XP:25:ARG:O	2.14	0.64
4:XD:9:CYS:SG	4:XD:31:CYS:O	2.54	0.64
31:YH:105:LEU:H	31:YH:105:LEU:CD1	2.09	0.64
36:YQ:23:GLY:HA3	36:YQ:101:ARG:NH1	2.12	0.64
50:Y4:18:CYS:SG	50:Y4:39:CYS:CB	2.85	0.64
1:QA:1392:G:H21	1:QA:1502:A:H8	1.43	0.64
25:RA:709:U:H3	25:RA:722:A:N6	1.94	0.64
28:RE:13:ARG:NH1	28:RE:21:VAL:HG12	2.11	0.64
34:RO:4:PRO:O	34:RO:5:GLN:HB2	1.96	0.64
3:XC:122:GLU:OE1	3:XC:126:ARG:NH2	2.29	0.64
5:XE:37:ARG:HA	5:XE:114:GLY:N	2.13	0.64
22:XV:3:G:HO2'	22:XV:4:G:H8	1.45	0.64
25:YA:1210:A:H5'	25:YA:1210:A:H8	1.62	0.64
27:YD:44:ASN:HB3	27:YD:49:ILE:HG22	1.78	0.64
30:YG:112:PRO:HB3	50:Y4:37:SER:HB2	1.79	0.64
4:QD:9:CYS:SG	4:QD:31:CYS:C	2.79	0.64
25:RA:910:A:N3	25:RA:2264:C:O2'	2.30	0.64
25:RA:1009:A:OP2	25:RA:1010:A:OP2	2.15	0.64
28:RE:35:GLN:CG	28:RE:37:ARG:HE	2.11	0.64
28:RE:50:GLY:HA3	28:RE:74:PRO:HG3	1.79	0.64
31:RH:148:ILE:O	31:RH:151:ILE:HG12	1.98	0.64
36:RQ:81:VAL:O	36:RQ:82:ARG:HG2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:990:A:OP2	25:YA:991:C:OP2	2.16	0.64
26:YB:77:U:P	45:YZ:19:ARG:HH22	2.20	0.64
30:YG:179:PRO:HG3	50:Y4:38:LYS:HZ2	1.61	0.64
36:YQ:81:VAL:O	36:YQ:82:ARG:HG2	1.97	0.64
50:Y4:18:CYS:HB3	50:Y4:39:CYS:CB	2.28	0.64
5:QE:78:HIS:CE1	5:QE:142:LEU:HD23	2.33	0.64
13:QM:122:LYS:O	13:QM:122:LYS:HD3	1.97	0.64
25:RA:273(C):C:H42	25:RA:363(C):G:H1	1.44	0.64
27:RD:65:ILE:HD11	27:RD:67:PHE:CE2	2.31	0.64
1:XA:1054:C:N3	24:XY:34:G:H1'	2.11	0.64
29:YF:175:THR:O	29:YF:176:LEU:HB2	1.96	0.64
35:YP:14:LYS:O	35:YP:16:ARG:N	2.31	0.64
1:QA:1321:C:H5"	1:QA:1322:C:H5"	1.80	0.64
10:QJ:53:PRO:HA	14:QN:42:ILE:HD12	1.79	0.64
27:RD:8:PRO:HB3	27:RD:14:ARG:HB2	1.79	0.64
54:R8:56:GLU:OE1	54:R8:56:GLU:N	2.30	0.64
2:XB:178:ARG:HG3	8:XH:72:PRO:HA	1.79	0.64
12:XL:126:LYS:C	12:XL:128:ALA:H	2.04	0.64
25:YA:2361:A:P	54:Y8:27:THR:HG23	2.38	0.64
28:YE:104:VAL:HG11	28:YE:188:VAL:CG2	2.27	0.64
29:YF:45:ARG:HH11	29:YF:45:ARG:CG	2.09	0.64
31:YH:117:PRO:HB3	31:YH:123:PHE:CE2	2.33	0.64
36:YQ:81:VAL:O	36:YQ:82:ARG:HD3	1.97	0.64
38:YS:78:LEU:HD11	38:YS:107:GLU:O	1.98	0.64
42:YW:41:LYS:HE3	51:Y5:25:LEU:HD21	1.80	0.64
1:QA:571:U:O4	1:QA:864:A:N6	2.31	0.64
27:RD:35:LYS:HG2	27:RD:64:ILE:N	2.12	0.64
36:RQ:10:ARG:O	36:RQ:11:LYS:HB2	1.98	0.64
40:RU:90:VAL:O	40:RU:92:ARG:N	2.30	0.64
1:XA:662:G:H2'	1:XA:663:A:H8	1.62	0.64
5:XE:43:LEU:HD22	5:XE:136:MET:HG3	1.80	0.64
15:XO:26:GLU:OE2	15:XO:77:ARG:NH1	2.31	0.64
28:YE:69:LYS:O	28:YE:71:GLY:N	2.27	0.64
29:YF:46:ARG:HG2	29:YF:46:ARG:NH1	2.00	0.64
29:YF:155:LEU:HD13	29:YF:174:VAL:CG1	2.27	0.64
36:YQ:80:GLU:C	36:YQ:81:VAL:HG13	2.22	0.64
36:YQ:86:GLY:C	36:YQ:88:GLY:N	2.52	0.64
5:QE:11:ILE:HD11	5:QE:31:LEU:HD12	1.80	0.64
10:QJ:77:PRO:O	10:QJ:79:ARG:NH1	2.30	0.64
25:RA:155:C:N3	25:RA:171:G:N2	2.42	0.64
36:RQ:23:GLY:HA3	36:RQ:101:ARG:NH1	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:YH:51:ARG:HG3	31:YH:51:ARG:HH11	1.62	0.64
31:YH:148:ILE:O	31:YH:151:ILE:HG12	1.97	0.64
1:QA:439:A:OP2	1:QA:493:G:N1	2.30	0.64
1:QA:838:G:C6	1:QA:842:C:H1'	2.33	0.64
1:QA:1128:C:OP1	9:QI:66:ARG:NH2	2.31	0.64
12:QL:18:VAL:HG23	12:QL:19:ARG:H	1.63	0.64
25:RA:2729:G:C1'	28:RE:187:ALA:HB2	2.22	0.64
8:XH:7:ALA:HB2	8:XH:85:ARG:HD3	1.80	0.64
25:YA:1550:C:H2'	25:YA:1551:C:H6	1.62	0.64
25:YA:2133:G:H1'	25:YA:2158:A:H61	1.63	0.64
31:YH:92:ILE:H	31:YH:92:ILE:HD12	1.63	0.64
37:YR:51:LEU:HD13	37:YR:66:VAL:HG13	1.79	0.64
40:YU:90:VAL:HG12	40:YU:91:ASP:H	1.61	0.64
45:YZ:94:GLU:HB2	45:YZ:130:PRO:HD2	1.80	0.64
50:Y4:18:CYS:HB3	50:Y4:39:CYS:HB3	1.78	0.64
1:QA:1227:A:OP1	19:QS:80:TYR:OH	2.06	0.64
25:RA:957:A:H5'	36:RQ:76:LYS:HD2	1.80	0.64
25:RA:2321:G:H2'	25:RA:2321:G:N3	2.10	0.64
31:RH:92:ILE:H	31:RH:92:ILE:HD12	1.63	0.64
3:XC:11:ARG:O	3:XC:13:GLY:N	2.30	0.64
29:YF:11:VAL:HG12	29:YF:12:LEU:N	2.13	0.64
31:YH:3:ARG:HA	31:YH:3:ARG:NE	2.12	0.64
36:YQ:86:GLY:C	36:YQ:88:GLY:H	2.03	0.64
40:YU:50:ARG:O	40:YU:54:LYS:NZ	2.31	0.64
2:QB:5:ILE:HG21	2:QB:221:LEU:HD23	1.78	0.64
36:RQ:30:GLY:HA3	36:RQ:106:VAL:O	1.98	0.64
37:RR:104:ARG:HD3	37:RR:109:ALA:HB3	1.79	0.64
3:XC:19:GLU:O	3:XC:40:ARG:NH2	2.30	0.64
28:YE:4:ILE:C	28:YE:5:LEU:HD23	2.22	0.64
28:YE:14:ILE:CG1	28:YE:15:PHE:H	2.08	0.64
27:RD:35:LYS:HG2	27:RD:64:ILE:H	1.63	0.63
28:RE:201:THR:HG21	28:RE:203:LYS:HB3	1.80	0.63
43:RX:43:VAL:HG13	43:RX:51:VAL:HG21	1.79	0.63
1:XA:27:G:H4'	4:XD:209:ARG:HG3	1.79	0.63
1:XA:80:G:O6	1:XA:89:U:O4	2.16	0.63
5:XE:147:ASP:O	5:XE:151:LEU:HG	1.97	0.63
25:YA:1332:G:H21	25:YA:1610:A:H8	1.44	0.63
25:YA:1796:U:H2'	25:YA:1797:C:C6	2.33	0.63
25:YA:2377:A:O2'	38:YS:111:GLU:O	2.13	0.63
25:YA:2420:C:H41	54:Y8:30:ARG:HD2	1.62	0.63
25:YA:2635:C:H5"	28:YE:78:LEU:HA	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:12:SER:C	27:YD:14:ARG:H	2.06	0.63
38:YS:26:LEU:HD22	38:YS:87:PHE:HD1	1.63	0.63
39:YT:36:GLU:HG3	39:YT:41:ARG:HE	1.62	0.63
40:YU:92:ARG:CZ	41:YV:11:GLN:H	2.11	0.63
54:Y8:48:PHE:N	54:Y8:48:PHE:CD1	2.66	0.63
38:RS:26:LEU:HB3	38:RS:87:PHE:HA	1.81	0.63
47:R1:29:GLY:O	47:R1:31:GLY:N	2.29	0.63
5:XE:78:HIS:HB3	8:XH:107:LEU:HD12	1.79	0.63
26:YB:52:A:H62	38:YS:33:LYS:HG3	1.63	0.63
27:YD:18:VAL:HG12	27:YD:19:ALA:O	1.98	0.63
27:YD:135:PHE:N	27:YD:135:PHE:HD1	1.96	0.63
27:YD:230:ASP:O	27:YD:231:HIS:HB2	1.98	0.63
28:YE:201:THR:HG21	28:YE:203:LYS:HB3	1.80	0.63
38:YS:42:ASP:C	38:YS:44:LYS:H	2.07	0.63
2:QB:24:TRP:H	2:QB:24:TRP:HD1	1.46	0.63
4:QD:9:CYS:HB2	4:QD:22:LYS:HZ1	1.63	0.63
10:QJ:4:ILE:HB	10:QJ:74:ILE:HG13	1.81	0.63
10:QJ:58:ASP:O	10:QJ:59:SER:CB	2.46	0.63
31:RH:3:ARG:NE	31:RH:3:ARG:HA	2.12	0.63
31:RH:117:PRO:HB3	31:RH:123:PHE:CE2	2.33	0.63
36:RQ:104:PHE:O	36:RQ:105:GLU:HB3	1.98	0.63
48:R2:65:ASN:HB3	48:R2:69:ARG:HH22	1.61	0.63
50:R4:71:ARG:HH11	50:R4:71:ARG:CG	1.97	0.63
25:YA:943:U:OP2	35:YP:36:LYS:NZ	2.31	0.63
28:YE:11:MET:HE3	28:YE:186:GLY:HA2	1.79	0.63
28:YE:50:GLY:HA3	28:YE:74:PRO:HG3	1.79	0.63
35:YP:5:ASP:O	35:YP:6:LEU:C	2.41	0.63
36:YQ:104:PHE:O	36:YQ:105:GLU:HB3	1.98	0.63
1:QA:1002:G:N2	1:QA:1038:C:N3	2.44	0.63
5:QE:102:ALA:HB1	5:QE:106:PRO:HG2	1.80	0.63
25:RA:2131:G:H4'	25:RA:2132:U:H4'	1.80	0.63
31:RH:153:LYS:HG3	31:RH:161:GLY:HA3	1.80	0.63
36:RQ:81:VAL:O	36:RQ:82:ARG:HD3	1.98	0.63
37:RR:33:ARG:HG3	37:RR:115:GLU:HB3	1.79	0.63
1:XA:1201:A:H4'	1:XA:1202:G:H5"	1.81	0.63
2:XB:79:ASP:HA	2:XB:82:ARG:HB2	1.80	0.63
2:XB:187:LEU:HA	2:XB:201:ILE:HB	1.79	0.63
12:XL:18:VAL:HG23	12:XL:19:ARG:H	1.62	0.63
13:XM:91:ARG:HB2	13:XM:98:VAL:HG13	1.80	0.63
25:YA:1094:U:O2'	25:YA:1096:A:OP1	2.15	0.63
25:YA:1833:U:O2'	25:YA:1969:A:N1	2.26	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:YF:67:GLN:O	29:YF:67:GLN:CG	2.32	0.63
33:YN:4:TYR:O	40:YU:64:ARG:NH1	2.32	0.63
50:Y4:48:ARG:O	50:Y4:50:VAL:N	2.31	0.63
1:QA:690:G:H2'	1:QA:691:G:O4'	1.99	0.63
1:QA:1422:G:H5"	34:RO:48:PRO:HB3	1.80	0.63
19:QS:40:ILE:HG23	19:QS:41:VAL:HG22	1.79	0.63
25:RA:1348:G:H2'	25:RA:1349:A:H5"	1.81	0.63
36:RQ:20:ALA:HB1	36:RQ:99:PRO:CB	2.28	0.63
36:RQ:81:VAL:HG23	46:R0:7:LEU:CD1	2.29	0.63
40:RU:66:ASN:O	40:RU:70:ARG:HB2	1.98	0.63
50:R4:63:TYR:C	50:R4:65:ASP:H	2.05	0.63
1:XA:949:A:N7	13:XM:106:ASN:ND2	2.46	0.63
1:XA:1525:G:OP1	11:XK:120:ARG:NH2	2.32	0.63
12:XL:62:SER:O	12:XL:64:TYR:HD1	1.82	0.63
25:YA:1138:G:N2	33:YN:106:MET:HE3	2.09	0.63
35:YP:101:VAL:HG23	35:YP:106:LEU:HB3	1.80	0.63
44:YY:91:GLU:HG3	44:YY:92:ASN:H	1.63	0.63
1:QA:973:G:H3'	1:QA:974:A:H5"	1.80	0.63
5:QE:71:LEU:HD11	5:QE:115:VAL:N	2.13	0.63
28:RE:131:ALA:HB1	28:RE:135:HIS:CE1	2.34	0.63
36:RQ:66:ILE:CG1	36:RQ:67:ARG:H	2.12	0.63
5:XE:37:ARG:HA	5:XE:114:GLY:H	1.62	0.63
25:YA:414:C:O2	25:YA:1864:U:O2'	2.15	0.63
25:YA:1820:U:C2	27:YD:202:LYS:HB3	2.34	0.63
27:YD:72:LYS:HG2	27:YD:103:ARG:NH2	2.13	0.63
31:YH:153:LYS:HG3	31:YH:161:GLY:HA3	1.80	0.63
50:Y4:23:GLU:O	50:Y4:25:TYR:N	2.31	0.63
12:QL:85:ILE:HD11	12:QL:98:TYR:HB2	1.81	0.63
50:R4:39:CYS:HB3	50:R4:41:PRO:HD2	1.79	0.63
54:R8:52:LYS:HG3	54:R8:52:LYS:O	1.97	0.63
1:XA:250:A:H4'	1:XA:251:G:O5'	1.99	0.63
12:XL:86:ARG:HB2	12:XL:101:VAL:CG2	2.28	0.63
13:XM:23:TYR:HB3	13:XM:67:GLU:HA	1.81	0.63
38:YS:88:ASP:CG	38:YS:90:GLY:H	2.06	0.63
40:YU:83:LEU:HD12	40:YU:113:ALA:HB2	1.79	0.63
40:YU:92:ARG:O	40:YU:94:ASN:N	2.29	0.63
1:QA:346:G:H1'	1:QA:347:G:H5'	1.80	0.63
8:QH:10:LEU:HD22	8:QH:83:ILE:HD11	1.80	0.63
12:QL:62:SER:O	12:QL:64:TYR:HD1	1.82	0.63
12:QL:86:ARG:HB2	12:QL:101:VAL:CG2	2.28	0.63
33:RN:13:TRP:HB2	33:RN:133:GLN:HG3	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:RN:133:GLN:HB2	33:RN:135:PRO:HD3	1.79	0.63
54:R8:48:PHE:N	54:R8:48:PHE:CD1	2.66	0.63
5:XE:69:VAL:O	5:XE:71:LEU:N	2.31	0.63
10:XJ:58:ASP:O	10:XJ:59:SER:CB	2.46	0.63
25:YA:451:C:H4'	29:YF:52:LYS:NZ	2.13	0.63
25:YA:910:A:C5	36:YQ:13:GLN:HG3	2.33	0.63
25:YA:1187:G:H5"	41:YV:81:TYR:CE1	2.34	0.63
25:YA:1354:A:OP1	27:YD:38:LYS:HE2	1.99	0.63
34:YO:13:ASN:ND2	34:YO:96:THR:O	2.30	0.63
36:YQ:20:ALA:HB1	36:YQ:99:PRO:CB	2.29	0.63
36:YQ:30:GLY:HA3	36:YQ:106:VAL:O	1.98	0.63
48:Y2:46:GLN:OE1	48:Y2:46:GLN:HA	1.97	0.63
13:QM:3:ARG:NH1	30:RG:113:ARG:NH2	2.47	0.63
36:RQ:20:ALA:HB1	36:RQ:99:PRO:CD	2.29	0.63
38:RS:15:ARG:HH11	38:RS:25:ARG:HH21	1.44	0.63
44:RY:51:VAL:HG13	44:RY:52:SER:H	1.64	0.63
25:YA:252:G:OP2	35:YP:50:ARG:NH1	2.32	0.63
25:YA:1043:C:H42	25:YA:1112:G:H1	1.47	0.63
26:YB:15:A:H5'	26:YB:16:G:C8	2.33	0.63
28:YE:35:GLN:CG	28:YE:37:ARG:NE	2.62	0.63
32:YI:130:TYR:HB3	32:YI:136:VAL:HG13	1.81	0.63
36:YQ:10:ARG:O	36:YQ:11:LYS:HB2	1.97	0.63
36:YQ:90:VAL:C	36:YQ:92:GLY:H	2.07	0.63
38:YS:22:GLY:O	38:YS:23:ARG:O	2.17	0.63
48:Y2:17:SER:HB2	48:Y2:18:PRO:CA	2.29	0.63
53:Y7:9:ARG:NH2	53:Y7:48:LYS:HD2	2.05	0.63
2:QB:24:TRP:H	2:QB:24:TRP:CD1	2.15	0.62
7:QG:155:ARG:HD3	7:QG:155:ARG:H	1.64	0.62
22:QV:75:C:H2'	22:QV:76:A:O4'	1.98	0.62
25:RA:622:G:OP2	35:RP:108:LYS:NZ	2.23	0.62
25:RA:883:G:H22	25:RA:892:G:H22	1.46	0.62
25:RA:2485:G:OP1	36:RQ:46:GLN:NE2	2.29	0.62
25:RA:2494:G:H2'	25:RA:2495:G:H8	1.62	0.62
1:XA:258:G:H2'	1:XA:259:G:H8	1.64	0.62
10:XJ:32:ALA:HB3	10:XJ:76:ASN:HB2	1.79	0.62
25:YA:67:U:H3	25:YA:74:A:H2	0.74	0.62
25:YA:1407:C:H42	25:YA:1595:G:H1	1.45	0.62
28:YE:13:ARG:HH12	28:YE:21:VAL:HG12	1.64	0.62
29:YF:129:PHE:O	29:YF:130:ALA:HB3	1.99	0.62
29:YF:132:VAL:HG23	29:YF:133:ASN:N	2.14	0.62
31:YH:86:GLU:O	31:YH:87:LEU:HB2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:Y6:41:PRO:HG2	52:Y6:45:LYS:H	1.63	0.62
4:QD:22:LYS:HG2	4:QD:26:CYS:SG	2.39	0.62
31:RH:136:ILE:H	31:RH:136:ILE:HD12	1.64	0.62
2:XB:212:GLN:NE2	2:XB:235:SER:HB2	2.14	0.62
25:YA:1509:C:H3'	25:YA:1510:A:H5"	1.80	0.62
25:YA:1728:G:H3'	25:YA:1729:A:H5"	1.80	0.62
28:YE:4:ILE:CD1	28:YE:28:ALA:HB1	2.29	0.62
35:YP:32:THR:O	35:YP:32:THR:OG1	2.12	0.62
54:Y8:59:LYS:HZ3	54:Y8:59:LYS:HB3	1.63	0.62
36:RQ:90:VAL:C	36:RQ:92:GLY:H	2.06	0.62
9:XI:24:GLY:N	9:XI:60:ASP:OD1	2.29	0.62
19:XS:13:ASP:OD1	19:XS:13:ASP:N	2.32	0.62
25:YA:128:C:H4'	53:Y7:49:ARG:HH12	1.63	0.62
25:YA:1434:A:H61	25:YA:1558:A:N6	1.97	0.62
25:YA:1818:U:H2'	27:YD:157:ARG:HG3	1.80	0.62
30:YG:3:LEU:HD12	30:YG:4:ASP:H	1.64	0.62
30:YG:115:ARG:NH2	30:YG:137:GLU:OE1	2.32	0.62
38:YS:48:LEU:HD12	38:YS:48:LEU:N	2.14	0.62
51:Y5:4:HIS:HB3	51:Y5:5:PRO:HD3	1.81	0.62
1:QA:7:G:H5'	1:QA:298:A:O4'	2.00	0.62
25:RA:1385:G:O2'	25:RA:1396:U:O2	2.12	0.62
29:RF:143:ALA:HB1	29:RF:148:LEU:HB2	1.81	0.62
35:RP:62:LEU:HD21	54:R8:25:MET:CB	2.26	0.62
1:XA:1014:A:H4'	19:XS:14:HIS:CD2	2.34	0.62
2:XB:12:GLU:O	2:XB:16:HIS:ND1	2.21	0.62
31:YH:136:ILE:H	31:YH:136:ILE:HD12	1.64	0.62
1:QA:877:C:H5"	8:QH:88:LYS:HD3	1.80	0.62
25:RA:1086:A:O2'	25:RA:1087:G:N7	2.31	0.62
25:RA:1628:G:H1	25:RA:1638:C:H42	1.47	0.62
25:RA:2111:C:N3	25:RA:2118:U:O2'	2.32	0.62
32:RI:52:ARG:HB2	32:RI:56:LYS:HB3	1.82	0.62
2:XB:92:TYR:CE1	2:XB:151:GLY:HA3	2.35	0.62
10:XJ:5:ARG:HH21	10:XJ:99:LYS:HD2	1.63	0.62
25:YA:2817:G:OP1	37:YR:99:LYS:NZ	2.27	0.62
35:YP:147:LEU:O	35:YP:148:LEU:HB2	1.97	0.62
39:YT:1:MET:O	39:YT:3:ARG:N	2.29	0.62
28:RE:35:GLN:CG	28:RE:37:ARG:NE	2.62	0.62
31:RH:137:ASP:HB3	31:RH:140:LYS:HB2	1.81	0.62
50:R4:38:LYS:C	50:R4:40:HIS:N	2.52	0.62
1:XA:7:G:H5'	1:XA:298:A:O4'	2.00	0.62
1:XA:1348:U:H3	1:XA:1374:A:H2	1.47	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:XG:111:ARG:NH1	7:XG:113:GLU:OE2	2.32	0.62
23:XX:19:A:O2'	23:XX:20:C:H5'	1.98	0.62
25:YA:2011:U:OP1	42:YW:42:ARG:NH1	2.33	0.62
27:YD:35:LYS:HA	27:YD:64:ILE:HG22	1.82	0.62
27:YD:133:LEU:HD21	27:YD:191:ALA:CB	2.29	0.62
28:YE:131:ALA:HB1	28:YE:135:HIS:CE1	2.34	0.62
48:Y2:41:ILE:HG12	48:Y2:44:LEU:HD12	1.82	0.62
48:Y2:69:ARG:HB2	48:Y2:69:ARG:CZ	2.29	0.62
8:QH:6:ILE:HB	8:QH:85:ARG:NH1	2.15	0.62
39:RT:54:ARG:HA	39:RT:59:THR:HG23	1.82	0.62
43:RX:53:LYS:HB2	43:RX:82:GLN:HB3	1.80	0.62
1:XA:1218:C:H2'	1:XA:1219:U:C6	2.34	0.62
2:XB:54:THR:HG21	2:XB:201:ILE:HD11	1.82	0.62
25:YA:273(C):C:H42	25:YA:363(C):G:H1	1.46	0.62
29:YF:28:ILE:HD13	29:YF:30:PRO:HD3	1.80	0.62
38:YS:100:ALA:HA	38:YS:103:GLU:CG	2.30	0.62
48:Y2:70:GLN:O	48:Y2:71:ASN:HB2	2.00	0.62
4:QD:3:ARG:HH11	4:QD:115:ARG:HD2	1.64	0.62
25:RA:2011:U:OP2	42:RW:16:LYS:NZ	2.33	0.62
28:RE:13:ARG:HH12	28:RE:21:VAL:HG12	1.64	0.62
5:XE:42:GLY:HA3	5:XE:66:MET:HG2	1.82	0.62
25:YA:958:U:OP2	36:YQ:14:ARG:NH1	2.33	0.62
25:YA:1678:G:N2	25:YA:1989:G:H22	1.97	0.62
1:QA:13:U:C4	1:QA:20:U:O4	2.53	0.62
1:QA:1352:C:OP1	21:QU:3:LYS:NZ	2.26	0.62
13:QM:3:ARG:HH22	30:RG:113:ARG:NH2	1.89	0.62
17:QQ:66:SER:O	17:QQ:70:ARG:NH1	2.33	0.62
25:RA:2298:A:H62	25:RA:2318:G:H8	1.46	0.62
1:XA:328:C:H4'	1:XA:329:A:H5'	1.82	0.62
1:XA:547:A:OP1	4:XD:73:ARG:NH2	2.32	0.62
1:XA:664:G:N2	1:XA:741:G:H1	1.96	0.62
25:YA:1057:A:N1	25:YA:1081:U:O4	2.33	0.62
25:YA:2583:G:N3	56:Z6:101:PPU:H2	2.15	0.62
28:YE:104:VAL:HG11	28:YE:188:VAL:HG23	1.82	0.62
25:RA:1026:U:H4'	25:RA:1027:A:OP1	2.00	0.62
27:RD:35:LYS:HD2	27:RD:104:TYR:CD1	2.35	0.62
28:RE:201:THR:HG22	28:RE:203:LYS:N	2.07	0.62
1:XA:45:U:H2'	1:XA:46:G:C8	2.35	0.62
27:YD:134:ARG:HD3	27:YD:135:PHE:CE1	2.35	0.62
27:YD:182:LEU:H	27:YD:272:ALA:HB3	1.62	0.62
29:YF:28:ILE:HG22	29:YF:112:MET:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:YP:65:ARG:NH2	54:Y8:46:ARG:HH12	1.98	0.62
36:YQ:63:LYS:HD2	45:YZ:175:VAL:HG21	1.82	0.62
46:Y0:27:GLU:HG3	46:Y0:68:GLU:HA	1.82	0.62
48:Y2:40:SER:C	48:Y2:42:GLY:N	2.51	0.62
1:QA:671:G:H1	1:QA:735:C:N4	1.95	0.61
1:QA:1292:U:OP1	7:QG:41:ARG:NH2	2.33	0.61
11:QK:121:PRO:HD2	11:QK:126:ARG:HD3	1.80	0.61
14:QN:13:THR:N	14:QN:14:PRO:HD2	2.15	0.61
22:QV:76:A:H2'	25:RA:2602:A:N6	2.15	0.61
25:RA:65:C:H1'	25:RA:456:C:H42	1.64	0.61
25:RA:67:U:N3	25:RA:74:A:H2	1.94	0.61
25:RA:587:C:OP2	35:RP:21:ARG:NH2	2.32	0.61
25:RA:1667:G:O2'	25:RA:1991:U:O4	2.18	0.61
28:RE:35:GLN:HG2	28:RE:37:ARG:NE	2.14	0.61
50:R4:23:GLU:O	50:R4:25:TYR:N	2.33	0.61
50:R4:61:ARG:O	50:R4:63:TYR:N	2.33	0.61
54:R8:33:ASN:O	54:R8:34:TRP:C	2.42	0.61
4:XD:111:ALA:HB2	4:XD:120:LEU:HD12	1.82	0.61
9:XI:29:ASN:OD1	9:XI:65:VAL:N	2.29	0.61
48:Y2:17:SER:CB	48:Y2:18:PRO:HA	2.29	0.61
50:Y4:37:SER:HB3	50:Y4:42:PHE:CD1	2.35	0.61
1:QA:1316:G:H22	1:QA:1319:A:H5"	1.65	0.61
4:QD:52:SER:H	4:QD:55:ALA:HB3	1.65	0.61
26:RB:50:G:H5"	38:RS:61:ASN:HD21	1.64	0.61
29:RF:107:LYS:HE3	29:RF:206:ILE:HD12	1.82	0.61
31:RH:152:ARG:O	31:RH:153:LYS:CD	2.48	0.61
1:XA:346:G:H1'	1:XA:347:G:H5'	1.82	0.61
27:YD:227:ASN:CB	27:YD:228:PRO:HD2	2.24	0.61
29:YF:32:LEU:CD1	29:YF:105:VAL:HG13	2.29	0.61
35:YP:71:VAL:HG13	35:YP:72:PRO:HD3	1.81	0.61
1:QA:1305:G:HO2'	1:QA:1306:A:H8	1.49	0.61
4:QD:108:LEU:HD21	4:QD:183:GLY:HA3	1.83	0.61
5:QE:101:ILE:CD1	5:QE:119:LEU:HD23	2.30	0.61
1:XA:1453:G:N7	20:XT:55:ILE:HD11	2.16	0.61
26:YB:38:C:H42	26:YB:44:G:H1	1.47	0.61
35:YP:65:ARG:NH2	54:Y8:46:ARG:NH1	2.48	0.61
54:Y8:33:ASN:O	54:Y8:34:TRP:C	2.42	0.61
1:QA:382:A:H2'	1:QA:383:A:C8	2.35	0.61
25:RA:574:C:N3	28:RE:145:LYS:NZ	2.47	0.61
31:RH:86:GLU:O	31:RH:87:LEU:HB2	1.99	0.61
31:RH:153:LYS:CB	31:RH:154:PRO:CD	2.69	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:486:U:H2'	1:XA:487:A:H8	1.65	0.61
1:XA:1299:A:H2'	1:XA:1301:U:H1'	1.81	0.61
25:YA:896:A:C8	45:YZ:146:ILE:HD12	2.35	0.61
27:YD:27:THR:O	27:YD:29:PRO:HD2	1.99	0.61
31:YH:137:ASP:HB3	31:YH:140:LYS:HB2	1.81	0.61
36:YQ:54:MET:O	36:YQ:57:HIS:HB3	2.00	0.61
10:QJ:40:LEU:HB2	10:QJ:69:ASN:HB3	1.82	0.61
25:RA:1711:C:H42	25:RA:1747:G:H1	1.48	0.61
25:RA:1798:U:H5'	27:RD:259:THR:HG22	1.81	0.61
47:R1:53:VAL:HG22	47:R1:74:VAL:HG13	1.83	0.61
55:R9:25:VAL:HB	55:R9:34:GLN:HB2	1.82	0.61
1:XA:280:C:C2	17:XQ:38:ARG:HG3	2.36	0.61
1:XA:1135:U:H4'	1:XA:1136:U:H5	1.66	0.61
25:YA:1021:A:H61	25:YA:1142(A):A:H61	1.46	0.61
25:YA:1196:C:HO2'	25:YA:1228:G:HO2'	1.39	0.61
25:YA:1204:A:H1'	25:YA:1206:G:C8	2.35	0.61
25:YA:2308:G:O6	25:YA:2311:A:N1	2.33	0.61
25:YA:2335:A:O2'	25:YA:2336:A:O5'	2.17	0.61
28:YE:174:ASP:CG	28:YE:175:VAL:N	2.58	0.61
31:YH:6:ARG:HG3	31:YH:7:LEU:N	2.15	0.61
36:YQ:66:ILE:CG1	36:YQ:67:ARG:H	2.12	0.61
48:Y2:41:ILE:HD11	48:Y2:44:LEU:HG	1.82	0.61
1:QA:1286:A:H8	1:QA:1287:A:H4'	1.65	0.61
14:QN:6:LEU:HD23	14:QN:23:ARG:HH22	1.64	0.61
25:RA:2817:G:O2'	25:RA:2836:U:O2	2.18	0.61
12:XL:85:ILE:HD11	12:XL:98:TYR:HB2	1.81	0.61
25:YA:503:A:H4'	25:YA:504:U:H5'	1.81	0.61
27:YD:25:THR:HG21	27:YD:81:ALA:CA	2.31	0.61
27:YD:70:TRP:CH2	27:YD:150:LYS:HA	2.35	0.61
27:YD:137:PRO:HB2	27:YD:140:THR:HG23	1.81	0.61
27:YD:147:LEU:CD1	27:YD:155:LEU:HD11	2.26	0.61
28:YE:69:LYS:C	28:YE:71:GLY:H	2.08	0.61
28:YE:95:ILE:HD12	28:YE:95:ILE:N	2.15	0.61
38:YS:17:ARG:HG3	38:YS:18:ILE:N	2.14	0.61
38:YS:49:VAL:HG22	38:YS:80:LEU:HD12	1.82	0.61
54:Y8:29:LYS:HD3	54:Y8:44:LYS:CB	2.30	0.61
3:QC:14:ILE:O	3:QC:16:ARG:N	2.33	0.61
11:QK:22:HIS:HB3	11:QK:29:ILE:HG23	1.83	0.61
21:QU:6:ARG:HE	21:QU:15:ARG:NH2	1.99	0.61
25:RA:807:U:H2'	25:RA:808:G:C8	2.34	0.61
31:RH:6:ARG:HG3	31:RH:7:LEU:N	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:R4:71:ARG:NH1	50:R4:71:ARG:CG	2.60	0.61
54:R8:29:LYS:HD3	54:R8:44:LYS:CB	2.31	0.61
2:XB:67:THR:HG21	2:XB:155:LEU:HG	1.81	0.61
8:XH:4:ASP:OD2	8:XH:85:ARG:NH1	2.33	0.61
15:XO:87:ILE:HG22	15:XO:88:ARG:H	1.64	0.61
20:XT:100:ILE:HG13	20:XT:102:GLY:N	2.16	0.61
25:YA:226:G:O2'	25:YA:228:A:N6	2.33	0.61
25:YA:2286:A:H8	52:Y6:37:ARG:HH11	1.48	0.61
27:YD:133:LEU:HD21	27:YD:191:ALA:HB2	1.82	0.61
28:YE:35:GLN:HG2	28:YE:37:ARG:NE	2.14	0.61
36:YQ:2:LEU:HD23	36:YQ:2:LEU:H	1.65	0.61
54:Y8:22:VAL:HG21	54:Y8:53:PRO:HB2	1.83	0.61
15:QO:82:ILE:O	15:QO:86:GLY:N	2.33	0.61
42:RW:25:ARG:NH2	42:RW:74:ALA:O	2.33	0.61
1:XA:1459:C:OP1	20:XT:27:LYS:NZ	2.34	0.61
27:YD:69:ARG:C	27:YD:71:ASP:H	2.08	0.61
28:YE:35:GLN:CG	28:YE:37:ARG:HE	2.11	0.61
28:YE:52:LEU:HB3	28:YE:54:GLN:OE1	2.00	0.61
29:YF:164:ARG:HG2	29:YF:164:ARG:HH11	1.66	0.61
31:YH:152:ARG:O	31:YH:153:LYS:CD	2.48	0.61
1:QA:1095:U:P	1:QA:1108:G:H1	2.24	0.61
1:QA:1175:G:H2'	1:QA:1176:A:C8	2.36	0.61
28:RE:69:LYS:C	28:RE:71:GLY:H	2.09	0.61
29:RF:101:LEU:O	29:RF:106:ARG:NH1	2.33	0.61
36:RQ:2:LEU:H	36:RQ:2:LEU:HD23	1.65	0.61
1:XA:132:C:H42	1:XA:230:G:H1	1.47	0.61
27:YD:2:ALA:HB3	27:YD:20:ASP:HB3	1.83	0.61
29:YF:119:ARG:HH11	29:YF:119:ARG:HG2	1.64	0.61
38:YS:89:ARG:O	38:YS:90:GLY:O	2.19	0.61
40:YU:52:ARG:HA	40:YU:55:ARG:HG3	1.83	0.61
1:QA:673:G:H2'	1:QA:674:G:C8	2.36	0.61
19:QS:5:LEU:CD1	50:R4:67:TYR:CD2	2.83	0.61
28:RE:52:LEU:HB3	28:RE:54:GLN:OE1	2.00	0.61
50:R4:64:GLY:C	50:R4:66:SER:H	2.08	0.61
54:R8:59:LYS:HZ3	54:R8:59:LYS:HB3	1.66	0.61
25:YA:273:G:H1	25:YA:364:C:H42	1.48	0.61
27:YD:54:ARG:HH11	27:YD:54:ARG:CG	2.14	0.61
28:YE:131:ALA:HB1	28:YE:135:HIS:HE1	1.65	0.61
47:Y1:83:GLU:O	47:Y1:85:LEU:N	2.34	0.61
3:QC:8:ILE:HG23	3:QC:16:ARG:HG2	1.83	0.60
4:QD:9:CYS:HB2	4:QD:22:LYS:NZ	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:QE:100:VAL:HB	5:QE:107:ARG:HH21	1.66	0.60
25:RA:1059:G:O6	25:RA:1079:C:N4	2.34	0.60
25:RA:2068:U:N3	25:RA:2430:A:H2	1.89	0.60
25:RA:2116:G:N1	25:RA:2162:G:OP1	2.31	0.60
25:RA:2286:A:H4'	25:RA:2287:A:O4'	2.00	0.60
36:RQ:80:GLU:CD	46:R0:7:LEU:HB3	2.26	0.60
38:RS:88:ASP:O	38:RS:89:ARG:HB3	2.01	0.60
40:RU:90:VAL:HG11	41:RV:40:LEU:HD12	1.82	0.60
36:YQ:20:ALA:HB1	36:YQ:99:PRO:CD	2.30	0.60
36:YQ:66:ILE:CG1	36:YQ:67:ARG:N	2.64	0.60
51:Y5:16:ARG:HG2	51:Y5:16:ARG:HH11	1.66	0.60
25:YA:2089:U:H6	25:YA:2089:U:H5"	1.66	0.60
25:YA:2439:A:C8	25:YA:2439:A:H5'	2.36	0.60
28:YE:63:LEU:CD1	28:YE:64:LYS:H	2.04	0.60
29:YF:108:LYS:HA	29:YF:108:LYS:HZ3	1.64	0.60
33:YN:8:GLN:O	33:YN:9:VAL:HG13	2.01	0.60
50:Y4:56:VAL:HA	50:Y4:60:GLN:HB2	1.84	0.60
1:QA:1023:G:H3'	1:QA:1024:G:H5"	1.82	0.60
1:QA:1118:C:H1'	1:QA:1179:A:C4	2.36	0.60
3:QC:11:ARG:O	3:QC:13:GLY:N	2.34	0.60
25:RA:2393:A:H5'	35:RP:62:LEU:HB3	1.82	0.60
36:RQ:66:ILE:CG1	36:RQ:67:ARG:N	2.64	0.60
37:RR:70:LEU:O	37:RR:72:ASP:N	2.31	0.60
1:XA:1297:C:O2'	7:XG:114:ARG:NH2	2.33	0.60
25:YA:221:A:H4'	25:YA:222:A:O5'	2.00	0.60
25:YA:2698:U:H2'	25:YA:2699:C:C6	2.36	0.60
27:YD:35:LYS:HG2	27:YD:64:ILE:CG2	2.31	0.60
28:YE:37:ARG:NE	28:YE:37:ARG:CA	2.64	0.60
38:YS:99:LYS:O	38:YS:102:ALA:N	2.34	0.60
28:RE:104:VAL:HG11	28:RE:188:VAL:HG23	1.82	0.60
31:RH:126:PRO:CD	31:RH:127:GLU:N	2.65	0.60
36:RQ:54:MET:O	36:RQ:57:HIS:HB3	2.00	0.60
45:RZ:111:VAL:HG13	45:RZ:112:ARG:N	2.16	0.60
1:XA:1008:C:N4	1:XA:1021:G:H1	1.98	0.60
8:XH:91:ARG:HB2	12:XL:7:ILE:HG13	1.83	0.60
25:YA:676:A:H8	25:YA:2069:G:H21	1.48	0.60
25:YA:2030:A:H4'	25:YA:2031:A:H8	1.66	0.60
27:YD:35:LYS:NZ	27:YD:65:ILE:HA	2.15	0.60
27:YD:263:ARG:CB	27:YD:263:ARG:HH11	2.15	0.60
28:YE:53:PRO:HG2	28:YE:54:GLN:NE2	2.17	0.60
31:YH:44:VAL:O	31:YH:44:VAL:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:1296:C:OP1	13:QM:44:ARG:NH2	2.35	0.60
28:RE:4:ILE:CD1	28:RE:28:ALA:HB1	2.29	0.60
31:RH:44:VAL:O	31:RH:44:VAL:HG22	2.01	0.60
39:RT:77:PRO:HB2	39:RT:80:SER:HB2	1.83	0.60
44:RY:87:LYS:O	44:RY:88:LYS:NZ	2.33	0.60
25:YA:1814:G:H4'	27:YD:51:VAL:HG21	1.82	0.60
25:YA:2815:C:H5'	51:Y5:29:THR:HG21	1.82	0.60
1:QA:1106:G:H5"	3:QC:172:ARG:HG2	1.83	0.60
9:QI:13:ALA:HB2	9:QI:68:GLY:HA3	1.82	0.60
15:QO:39:LEU:HD13	15:QO:56:LEU:HB2	1.83	0.60
25:RA:530:G:N1	25:RA:2022:U:OP1	2.35	0.60
25:RA:1796:U:H2'	25:RA:1797:C:C6	2.36	0.60
27:RD:44:ASN:HB3	27:RD:49:ILE:HA	1.83	0.60
30:RG:114:ILE:HD13	30:RG:140:ILE:HG21	1.82	0.60
35:RP:85:LEU:HA	35:RP:88:LEU:HD22	1.83	0.60
17:XQ:11:VAL:HG12	17:XQ:85:VAL:HG13	1.84	0.60
25:YA:1257:C:O2'	29:YF:84:VAL:HG12	2.02	0.60
25:YA:1364:G:N7	47:Y1:2:SER:N	2.50	0.60
25:YA:2729:G:C1'	28:YE:187:ALA:HB2	2.32	0.60
27:YD:35:LYS:NZ	27:YD:64:ILE:O	2.32	0.60
29:YF:175:THR:O	29:YF:176:LEU:CB	2.49	0.60
38:YS:11:LYS:HB2	38:YS:91:PRO:HD3	1.84	0.60
1:QA:191:G:O2'	20:QT:101:GLY:O	2.20	0.60
1:QA:666:G:H5'	1:QA:726:C:H1'	1.84	0.60
1:QA:1104:G:H4'	2:QB:111:ARG:NH1	2.16	0.60
2:QB:15:VAL:H	2:QB:16:HIS:CE1	2.20	0.60
12:QL:54:LYS:CD	12:QL:54:LYS:N	2.65	0.60
13:QM:8:GLU:OE1	30:RG:115:ARG:HD3	2.01	0.60
25:RA:2419:U:O4	54:R8:30:ARG:NE	2.35	0.60
29:RF:9:ILE:HD11	29:RF:125:LEU:HG	1.83	0.60
1:XA:201:C:N4	1:XA:216:G:H1	1.96	0.60
12:XL:5:PRO:HA	12:XL:9:GLN:NE2	2.17	0.60
27:YD:25:THR:HG21	27:YD:81:ALA:HA	1.84	0.60
27:YD:72:LYS:HE3	27:YD:75:ILE:HD12	1.82	0.60
27:YD:166:GLN:HE21	27:YD:166:GLN:CA	2.14	0.60
39:YT:84:GLN:OE1	39:YT:85:LYS:NZ	2.34	0.60
1:QA:792:A:H4'	1:QA:793:U:O5'	2.00	0.60
1:QA:1004:A:P	1:QA:1025:U:H3	2.24	0.60
1:QA:1086:U:H3	1:QA:1099:G:H22	1.49	0.60
7:QG:26:PHE:O	7:QG:30:ILE:HG12	2.01	0.60
10:QJ:42:THR:HG23	10:QJ:68:HIS:HA	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2438:U:O3'	25:RA:2439:A:H3'	2.01	0.60
28:RE:131:ALA:HB1	28:RE:135:HIS:HE1	1.65	0.60
1:XA:987:G:H1	1:XA:1218:C:H42	1.49	0.60
5:XE:41:VAL:CG2	5:XE:113:ALA:HB2	2.31	0.60
25:YA:2415:G:H4'	35:YP:67:MET:N	2.17	0.60
25:YA:2593:U:H2'	25:YA:2594:C:H6	1.67	0.60
27:YD:35:LYS:HD3	27:YD:63:ARG:CB	2.32	0.60
27:YD:147:LEU:HD13	27:YD:155:LEU:CD1	2.29	0.60
28:YE:93:VAL:N	28:YE:95:ILE:CD1	2.65	0.60
28:YE:93:VAL:N	28:YE:95:ILE:HD12	2.17	0.60
30:YG:28:VAL:HG23	30:YG:29:TRP:CD1	2.36	0.60
34:YO:96:THR:O	34:YO:97:ARG:HB3	2.01	0.60
36:YQ:63:LYS:HE2	36:YQ:65:PHE:CE1	2.37	0.60
45:YZ:144:LEU:HD11	45:YZ:149:SER:HA	1.83	0.60
54:Y8:53:PRO:CD	54:Y8:54:GLU:H	2.15	0.60
1:QA:1446:A:H4'	39:RT:125:ARG:HH22	1.67	0.60
1:QA:1446:A:O2'	1:QA:1447:G:O5'	2.20	0.60
25:RA:2107:C:N3	25:RA:2182:G:N2	2.40	0.60
28:RE:53:PRO:HG2	28:RE:54:GLN:NE2	2.16	0.60
31:RH:89:ILE:O	31:RH:91:GLY:N	2.35	0.60
32:RI:13:GLY:HA3	32:RI:17:GLN:CD	2.26	0.60
36:RQ:97:VAL:HG21	36:RQ:103:MET:HE3	1.83	0.60
1:XA:582:U:H2'	1:XA:583:A:C8	2.36	0.60
1:XA:1226:C:O2'	13:XM:111:LYS:NZ	2.34	0.60
5:XE:43:LEU:HD22	5:XE:136:MET:HG2	1.84	0.60
12:XL:54:LYS:CD	12:XL:54:LYS:N	2.64	0.60
25:YA:2126:A:H4'	25:YA:2127:G:O5'	2.01	0.60
27:YD:21:PHE:HB3	27:YD:24:ILE:HG13	1.83	0.60
28:YE:68:ALA:O	28:YE:69:LYS:HG3	2.02	0.60
55:Y9:35:ARG:HH21	55:Y9:37:GLY:HA3	1.67	0.60
25:RA:181:A:H1'	25:RA:435:C:H5'	1.83	0.60
25:RA:2636:U:OP1	28:RE:79:ARG:HA	2.01	0.60
27:RD:70:TRP:CH2	27:RD:150:LYS:HA	2.36	0.60
28:RE:26:ILE:HD13	28:RE:26:ILE:C	2.26	0.60
28:RE:37:ARG:NE	28:RE:37:ARG:CA	2.64	0.60
28:RE:63:LEU:CD1	28:RE:64:LYS:H	2.04	0.60
29:RF:184:TYR:CE2	29:RF:188:ARG:HD2	2.37	0.60
31:RH:30:LYS:CD	31:RH:81:GLU:H	2.15	0.60
31:RH:117:PRO:HB3	31:RH:123:PHE:CD2	2.37	0.60
2:XB:74:LYS:HE3	2:XB:166:ASP:HB2	1.82	0.60
2:XB:235:SER:O	2:XB:237:ALA:N	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:XM:3:ARG:HA	13:XM:9:ILE:HG21	1.83	0.60
25:YA:2275:C:O2	36:YQ:83:MET:HG3	2.02	0.60
30:YG:6:ALA:H	50:Y4:23:GLU:CG	2.13	0.60
37:YR:42:LYS:HA	37:YR:45:ARG:HD2	1.84	0.60
38:YS:59:LYS:HG2	38:YS:60:GLY:N	2.13	0.60
28:RE:68:ALA:O	28:RE:69:LYS:HG3	2.02	0.59
28:RE:69:LYS:O	28:RE:71:GLY:N	2.27	0.59
28:RE:95:ILE:HD12	28:RE:95:ILE:N	2.15	0.59
28:RE:116:VAL:O	28:RE:117:MET:CB	2.49	0.59
1:XA:388:G:O2'	1:XA:389:A:OP2	2.17	0.59
25:YA:27:G:H22	25:YA:512:G:H2'	1.67	0.59
27:YD:35:LYS:CG	27:YD:64:ILE:N	2.56	0.59
29:YF:63:LYS:HE2	29:YF:67:GLN:HB3	1.83	0.59
31:YH:30:LYS:CD	31:YH:81:GLU:H	2.15	0.59
45:YZ:80:ARG:HH21	45:YZ:82:ARG:HH22	1.50	0.59
54:Y8:22:VAL:CG2	54:Y8:53:PRO:HB2	2.32	0.59
25:RA:1206:G:H1	25:RA:1240:U:H3	1.48	0.59
25:RA:2667:C:O2	31:RH:109:PHE:HB3	2.01	0.59
30:RG:64:THR:HG23	30:RG:66:GLN:H	1.66	0.59
32:RI:115:ALA:HB3	32:RI:128:LEU:HD12	1.84	0.59
36:RQ:81:VAL:HG23	36:RQ:82:ARG:H	1.67	0.59
39:RT:105:LEU:O	39:RT:107:ASP:N	2.36	0.59
1:XA:1336:C:H1'	1:XA:1337:G:C2	2.36	0.59
25:YA:2335:A:O2'	25:YA:2336:A:H2'	2.01	0.59
27:YD:172:TYR:CD1	27:YD:186:HIS:HA	2.37	0.59
28:YE:51:PHE:HD2	28:YE:52:LEU:HG	1.67	0.59
29:YF:123:LEU:HD12	29:YF:124:LEU:N	2.17	0.59
38:YS:110:LEU:HA	38:YS:112:PHE:CE2	2.37	0.59
25:RA:27:G:N2	25:RA:513:A:OP2	2.36	0.59
28:RE:61:ARG:HB2	28:RE:62:PRO:CD	2.33	0.59
31:RH:4:ILE:HD13	31:RH:4:ILE:N	2.18	0.59
36:RQ:63:LYS:HE2	36:RQ:65:PHE:CE1	2.37	0.59
45:RZ:5:LEU:HD11	45:RZ:39:VAL:HB	1.83	0.59
50:R4:35:VAL:O	50:R4:37:SER:N	2.26	0.59
1:XA:1015:A:H2'	1:XA:1016:A:C8	2.37	0.59
25:YA:49:A:N7	25:YA:120:U:C4	2.70	0.59
25:YA:443:A:C5	29:YF:45:ARG:HD2	2.38	0.59
25:YA:1265:A:H8	25:YA:1265:A:OP1	1.85	0.59
26:YB:40:U:O2'	26:YB:45:A:N6	2.33	0.59
28:YE:26:ILE:HD13	28:YE:26:ILE:C	2.26	0.59
28:YE:116:VAL:O	28:YE:117:MET:CB	2.49	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:YF:11:VAL:HG11	29:YF:18:ARG:HE	1.67	0.59
32:YI:129:THR:HA	32:YI:137:PRO:HA	1.84	0.59
42:YW:86:LEU:HD12	42:YW:87:PRO:HD2	1.83	0.59
1:QA:618:C:H5'	1:QA:619:U:H5"	1.84	0.59
3:QC:50:ALA:HB2	3:QC:75:VAL:HB	1.85	0.59
25:RA:747:U:C2	51:R5:2:ALA:N	2.70	0.59
25:RA:1649:G:O2'	37:RR:107:ASP:OD2	2.19	0.59
25:RA:2350:C:C5	54:R8:42:ARG:NH1	2.70	0.59
31:RH:82:GLY:O	31:RH:135:GLY:O	2.20	0.59
1:XA:582:U:H2'	1:XA:583:A:H8	1.68	0.59
6:XF:61:LEU:HB3	6:XF:63:TYR:HE1	1.66	0.59
12:XL:70:ILE:HD13	12:XL:77:LEU:HD12	1.83	0.59
25:YA:630:G:N2	25:YA:633:A:OP2	2.32	0.59
25:YA:2250:G:C5	36:YQ:82:ARG:HD2	2.37	0.59
31:YH:4:ILE:HD13	31:YH:4:ILE:N	2.18	0.59
31:YH:55:PRO:HG2	31:YH:61:HIS:CE1	2.37	0.59
35:YP:5:ASP:O	35:YP:6:LEU:O	2.19	0.59
35:YP:92:GLU:HA	35:YP:123:LEU:HD23	1.84	0.59
54:Y8:56:GLU:O	54:Y8:59:LYS:N	2.35	0.59
9:QI:9:ARG:HB3	9:QI:14:VAL:HG13	1.84	0.59
25:RA:602:G:O2'	25:RA:604:G:O2'	2.20	0.59
32:RI:92:VAL:HG13	32:RI:120:ILE:HG23	1.84	0.59
36:RQ:80:GLU:OE1	46:R0:6:GLY:O	2.20	0.59
38:RS:38:GLN:OE1	38:RS:47:THR:OG1	2.18	0.59
51:R5:40:LYS:NZ	51:R5:46:CYS:HB3	2.18	0.59
51:R5:52:TYR:O	51:R5:53:ALA:HB3	2.01	0.59
2:XB:92:TYR:HE1	2:XB:151:GLY:HA3	1.66	0.59
25:YA:521:G:H2'	25:YA:522:G:H8	1.68	0.59
25:YA:1020:A:N1	25:YA:1141:U:H2'	2.17	0.59
25:YA:2154:G:H2'	25:YA:2155:G:H8	1.67	0.59
25:YA:2185:C:H2'	25:YA:2186:G:C8	2.37	0.59
25:YA:2712:U:O2'	25:YA:2712(A):A:H8	1.85	0.59
31:YH:86:GLU:O	31:YH:131:VAL:O	2.20	0.59
31:YH:124:GLU:HB3	31:YH:132:ARG:HG3	1.84	0.59
53:Y7:9:ARG:HH21	53:Y7:48:LYS:HB2	1.62	0.59
1:QA:1157:A:H62	1:QA:1178:G:H21	1.49	0.59
1:QA:1222:G:OP1	19:QS:77:THR:HG21	2.02	0.59
1:QA:1384:C:H2'	1:QA:1385:G:H8	1.67	0.59
25:RA:307:G:H21	25:RA:330:A:H62	1.49	0.59
25:RA:442:G:H1'	29:RF:48:THR:HG21	1.85	0.59
25:RA:2150:U:H2'	25:RA:2151:G:H8	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2636:U:OP1	28:RE:79:ARG:HG3	2.03	0.59
1:XA:597:G:H1	1:XA:643:C:H42	1.49	0.59
25:YA:323:G:H2'	29:YF:169:ASN:OD1	2.02	0.59
25:YA:2343:C:O2'	25:YA:2373:G:O2'	2.21	0.59
31:YH:159:GLU:O	31:YH:160:LYS:HG2	2.03	0.59
49:Y3:6:VAL:HG13	49:Y3:56:VAL:HG13	1.84	0.59
50:Y4:18:CYS:CB	50:Y4:39:CYS:CB	2.79	0.59
25:RA:1071:G:N1	25:RA:1091:G:N7	2.51	0.59
25:RA:1416:G:H2'	25:RA:1417:C:C6	2.38	0.59
25:RA:2126:A:H4'	25:RA:2127:G:O5'	2.01	0.59
25:RA:2128:C:N4	25:RA:2160:G:H1	2.00	0.59
25:RA:2308:G:O6	25:RA:2311:A:N1	2.35	0.59
25:RA:2832:U:H4'	25:RA:2833:G:H5"	1.84	0.59
28:RE:36:ARG:H	28:RE:37:ARG:HH21	1.50	0.59
28:RE:116:VAL:CG2	28:RE:122:PHE:CD2	2.86	0.59
52:R6:52:VAL:HG22	52:R6:53:LYS:HG3	1.84	0.59
54:R8:22:VAL:CG2	54:R8:53:PRO:HB2	2.32	0.59
54:R8:22:VAL:HG21	54:R8:53:PRO:HB2	1.83	0.59
1:XA:564:C:O2'	8:XH:91:ARG:NH2	2.36	0.59
5:XE:45:PHE:CE2	5:XE:47:LYS:HD2	2.38	0.59
25:YA:481:G:OP2	44:YY:47:LYS:HG3	2.02	0.59
27:YD:174:ILE:HD12	27:YD:174:ILE:N	2.16	0.59
27:YD:177:LEU:HD11	27:YD:183:ARG:HB2	1.85	0.59
42:YW:111:HIS:CD2	42:YW:112:GLY:H	2.20	0.59
53:Y7:35:ARG:HG3	53:Y7:42:LEU:HD11	1.85	0.59
1:QA:266:G:O2'	1:QA:267:C:OP2	2.20	0.59
1:QA:501:C:H2'	1:QA:502:G:C8	2.35	0.59
12:QL:70:ILE:HD13	12:QL:77:LEU:HD12	1.83	0.59
25:RA:2518:A:H4'	25:RA:2519:U:OP1	2.01	0.59
26:RB:48:A:H2'	26:RB:49:C:C6	2.37	0.59
28:RE:93:VAL:N	28:RE:95:ILE:CD1	2.65	0.59
28:RE:174:ASP:CG	28:RE:175:VAL:N	2.58	0.59
31:RH:55:PRO:HG2	31:RH:61:HIS:CE1	2.37	0.59
42:RW:86:LEU:HD12	42:RW:87:PRO:HD2	1.85	0.59
50:R4:65:ASP:O	50:R4:66:SER:CB	2.51	0.59
25:YA:309:G:N3	25:YA:329:G:O2'	2.35	0.59
25:YA:442:G:O4'	29:YF:46:ARG:HD3	2.03	0.59
27:YD:27:THR:CG2	27:YD:83:GLU:HB3	2.33	0.59
31:YH:126:PRO:CD	31:YH:127:GLU:N	2.64	0.59
36:YQ:97:VAL:HG21	36:YQ:103:MET:HE3	1.83	0.59
38:YS:88:ASP:O	38:YS:89:ARG:CB	2.48	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:Y2:17:SER:HB2	48:Y2:18:PRO:HA	1.85	0.59
1:QA:1322:C:O2'	1:QA:1323:G:H5'	2.03	0.59
22:QV:53:G:H4'	22:QV:54:U:OP1	2.03	0.59
25:RA:900:A:H3'	25:RA:901:A:H8	1.66	0.59
25:RA:1210:A:H5"	25:RA:1210:A:C8	2.38	0.59
25:RA:2119:A:N6	25:RA:2170:A:N7	2.51	0.59
25:YA:1169:G:N2	25:YA:1180:C:N3	2.44	0.59
25:YA:2361:A:OP1	54:Y8:27:THR:HG23	2.03	0.59
25:YA:2599:G:OP2	27:YD:236:GLY:HA2	2.02	0.59
27:YD:137:PRO:HB2	27:YD:140:THR:CG2	2.33	0.59
28:YE:36:ARG:H	28:YE:37:ARG:HH21	1.49	0.59
28:YE:61:ARG:HB2	28:YE:62:PRO:CD	2.33	0.59
31:YH:82:GLY:O	31:YH:135:GLY:O	2.20	0.59
36:YQ:81:VAL:HG23	36:YQ:82:ARG:H	1.67	0.59
48:Y2:64:LEU:CD2	48:Y2:68:ARG:HD2	2.33	0.59
12:QL:54:LYS:N	12:QL:54:LYS:HD2	2.18	0.59
13:QM:49:THR:HB	13:QM:52:GLU:HG3	1.85	0.59
25:RA:242:G:C5'	54:R8:62:LEU:CD2	2.81	0.59
25:RA:1244:G:H4'	35:RP:7:ARG:HB2	1.84	0.59
25:RA:2733:A:N1	28:RE:203:LYS:O	2.36	0.59
27:RD:85:ASP:HB2	27:RD:92:ILE:HD13	1.84	0.59
31:RH:92:ILE:HG22	31:RH:93:GLY:N	2.18	0.59
31:RH:127:GLU:HG2	31:RH:128:PRO:CG	2.33	0.59
47:R1:92:LYS:HG3	47:R1:96:LYS:HB2	1.84	0.59
50:R4:12:ALA:CB	50:R4:29:PRO:HA	2.33	0.59
53:R7:22:MET:HE2	53:R7:31:LEU:HD12	1.85	0.59
10:XJ:76:ASN:O	10:XJ:78:ASN:ND2	2.36	0.59
13:XM:13:LYS:HA	13:XM:44:ARG:HD2	1.83	0.59
29:YF:63:LYS:HE2	29:YF:67:GLN:CB	2.32	0.59
31:YH:4:ILE:HG13	31:YH:6:ARG:CD	2.33	0.59
33:YN:40:PRO:HB3	40:YU:68:ALA:HB2	1.85	0.59
1:QA:1004:A:H1'	1:QA:1036:G:H22	1.68	0.58
5:QE:107:ARG:CZ	5:QE:107:ARG:HB2	2.30	0.58
7:QG:73:MET:HG2	7:QG:90:GLU:HA	1.83	0.58
19:QS:68:GLY:CA	50:R4:68:ARG:CG	2.74	0.58
25:RA:74:A:H4'	25:RA:75:G:O5'	2.03	0.58
28:RE:93:VAL:N	28:RE:95:ILE:HD12	2.17	0.58
1:XA:527:G:O6	12:XL:49:ASN:ND2	2.32	0.58
14:XN:23:ARG:HD2	14:XN:28:GLY:O	2.03	0.58
25:YA:2111:C:N3	25:YA:2118:U:O2'	2.36	0.58
27:YD:27:THR:CG2	27:YD:28:GLU:N	2.66	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:134:ILE:C	28:YE:134:ILE:HD12	2.28	0.58
30:YG:136:ARG:O	30:YG:154:GLY:HA2	2.02	0.58
31:YH:89:ILE:O	31:YH:91:GLY:N	2.35	0.58
48:Y2:32:LEU:HD11	48:Y2:54:LYS:HG3	1.84	0.58
25:RA:1470:G:N2	25:RA:1522:G:OP2	2.36	0.58
25:RA:2439:A:H5'	25:RA:2439:A:C8	2.37	0.58
27:RD:44:ASN:CB	27:RD:49:ILE:HA	2.33	0.58
27:RD:71:ASP:OD2	27:RD:103:ARG:NH2	2.36	0.58
28:RE:72:VAL:O	28:RE:73:GLU:O	2.21	0.58
31:RH:86:GLU:O	31:RH:131:VAL:O	2.21	0.58
35:RP:14:LYS:O	35:RP:16:ARG:N	2.36	0.58
50:R4:22:ILE:HG22	50:R4:23:GLU:N	2.18	0.58
50:R4:48:ARG:NH1	50:R4:52:THR:H	2.01	0.58
12:XL:18:VAL:O	12:XL:19:ARG:HB2	2.03	0.58
25:YA:2726:U:H4'	34:YO:1:MET:HE1	1.85	0.58
27:YD:165:ILE:HA	27:YD:175:LEU:HD23	1.83	0.58
29:YF:174:VAL:HG13	29:YF:174:VAL:O	2.04	0.58
12:QL:5:PRO:HA	12:QL:9:GLN:NE2	2.17	0.58
16:QP:21:VAL:O	16:QP:33:ILE:HG12	2.02	0.58
25:RA:1264:G:H5'	51:R5:11:THR:CG2	2.32	0.58
25:RA:2293:C:H5"	38:RS:89:ARG:HH12	1.69	0.58
45:RZ:45:ASP:OD1	45:RZ:49:ARG:NE	2.36	0.58
1:XA:93:U:H2'	1:XA:95:G:O4'	2.03	0.58
27:YD:44:ASN:HB3	27:YD:49:ILE:CA	2.27	0.58
31:YH:92:ILE:HG22	31:YH:93:GLY:N	2.18	0.58
37:YR:67:LEU:HD13	37:YR:76:VAL:HG21	1.85	0.58
41:YV:44:LYS:O	41:YV:46:VAL:N	2.36	0.58
43:YX:61:GLY:N	43:YX:75:ASP:OD1	2.36	0.58
47:Y1:51:VAL:HG11	47:Y1:74:VAL:HG21	1.84	0.58
28:RE:6:GLY:HA3	28:RE:26:ILE:HD11	1.85	0.58
30:RG:136:ARG:O	30:RG:154:GLY:HA2	2.03	0.58
50:R4:63:TYR:C	50:R4:65:ASP:N	2.62	0.58
1:XA:662:G:H2'	1:XA:663:A:C8	2.38	0.58
12:XL:45:PRO:HD3	12:XL:51:ALA:O	2.03	0.58
20:XT:10:LEU:HG	20:XT:12:ALA:H	1.68	0.58
25:YA:307:G:H21	25:YA:330:A:N6	2.01	0.58
25:YA:2784:C:H5"	28:YE:41:LYS:NZ	2.18	0.58
31:YH:117:PRO:HB3	31:YH:123:PHE:CD2	2.37	0.58
50:R4:42:PHE:CG	50:R4:43:TYR:N	2.72	0.58
51:R5:60:VAL:OXT	51:R5:60:VAL:HG13	2.03	0.58
54:R8:56:GLU:O	54:R8:59:LYS:N	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:486:U:H2'	1:XA:487:A:C8	2.39	0.58
2:XB:15:VAL:H	2:XB:16:HIS:CE1	2.22	0.58
2:XB:72:GLY:HA2	2:XB:165:VAL:HG22	1.85	0.58
8:XH:39:LEU:HB3	8:XH:45:ILE:HG12	1.85	0.58
12:XL:82:VAL:HG23	12:XL:106:ASP:OD2	2.04	0.58
25:YA:1266:G:O5'	42:YW:15:ARG:NH2	2.36	0.58
26:YB:81:G:O6	26:YB:95:U:O2	2.21	0.58
36:YQ:66:ILE:HA	36:YQ:104:PHE:HA	1.85	0.58
48:Y2:51:ARG:HA	48:Y2:54:LYS:HB2	1.86	0.58
48:Y2:69:ARG:NH1	48:Y2:69:ARG:CB	2.67	0.58
50:Y4:42:PHE:O	50:Y4:44:THR:N	2.36	0.58
1:QA:1348:U:H4'	9:QI:120:ARG:HD2	1.84	0.58
2:QB:77:ALA:HB2	2:QB:211:ILE:HD13	1.86	0.58
4:QD:105:VAL:HG13	4:QD:110:PHE:HB2	1.84	0.58
12:QL:33:ARG:O	12:QL:85:ILE:HG22	2.03	0.58
25:RA:1378:A:OP1	53:R7:10:ARG:NH2	2.36	0.58
26:RB:8:U:O3'	38:RS:25:ARG:NH2	2.32	0.58
27:RD:35:LYS:HD2	27:RD:104:TYR:CE1	2.39	0.58
27:RD:182:LEU:N	27:RD:272:ALA:HB3	2.17	0.58
31:RH:159:GLU:O	31:RH:160:LYS:HG2	2.03	0.58
32:RI:52:ARG:O	32:RI:56:LYS:N	2.37	0.58
35:RP:47:ASP:OD2	35:RP:50:ARG:NH2	2.36	0.58
36:RQ:81:VAL:C	36:RQ:82:ARG:CG	2.76	0.58
54:R8:46:ARG:O	54:R8:47:LYS:HB3	2.03	0.58
3:XC:14:ILE:HG12	3:XC:15:THR:H	1.68	0.58
6:XF:36:ARG:NH1	6:XF:38:GLU:OE2	2.36	0.58
9:XI:43:ALA:HA	9:XI:74:ILE:HD13	1.85	0.58
25:YA:771:G:OP1	53:Y7:14:LYS:HE3	2.04	0.58
25:YA:1490:A:O2'	27:YD:99:ASP:OD1	2.20	0.58
28:YE:51:PHE:CD2	28:YE:52:LEU:HG	2.38	0.58
29:YF:34:TRP:CZ3	35:YP:8:PRO:HB3	2.39	0.58
31:YH:85:LYS:HA	31:YH:86:GLU:OE1	2.04	0.58
36:YQ:55:VAL:HG22	36:YQ:56:ARG:N	2.18	0.58
10:QJ:49:VAL:HG13	14:QN:41:ARG:HB2	1.86	0.58
25:RA:137(A):G:H1'	43:RX:41:ASN:ND2	2.19	0.58
25:RA:2313:C:H5"	30:RG:91:ARG:HD3	1.85	0.58
25:RA:2667:C:H1'	31:RH:109:PHE:CD2	2.32	0.58
31:RH:4:ILE:HG13	31:RH:6:ARG:CD	2.33	0.58
31:RH:125:VAL:HG12	31:RH:126:PRO:CG	2.34	0.58
20:XT:63:ILE:HG22	20:XT:77:ALA:HB1	1.86	0.58
25:YA:2068:U:H3	25:YA:2430:A:H2	1.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2469:A:H5"	25:YA:2470:G:C8	2.39	0.58
27:YD:34:VAL:C	27:YD:35:LYS:HG3	2.28	0.58
27:YD:242:ARG:N	27:YD:242:ARG:HD2	2.18	0.58
29:YF:89:VAL:HG12	29:YF:90:PHE:N	2.18	0.58
29:YF:179:GLU:H	29:YF:179:GLU:CD	2.11	0.58
32:YI:129:THR:HG22	32:YI:137:PRO:HB3	1.86	0.58
36:YQ:47:ILE:CD1	36:YQ:70:PRO:HD3	2.34	0.58
37:YR:27:SER:HB3	37:YR:34:ILE:HD11	1.84	0.58
45:RZ:150:LEU:HD21	45:RZ:172:ALA:HB3	1.86	0.58
51:R5:50:GLY:O	51:R5:51:TYR:HB2	2.03	0.58
1:XA:1221:G:O3'	19:XS:77:THR:HG21	2.04	0.58
1:XA:1414:U:H2'	1:XA:1415:G:H8	1.69	0.58
2:XB:96:ARG:H	2:XB:96:ARG:HD2	1.67	0.58
3:XC:95:THR:HG22	3:XC:97:LYS:HG3	1.84	0.58
12:XL:54:LYS:N	12:XL:54:LYS:HD2	2.18	0.58
20:XT:56:MET:HG3	20:XT:88:VAL:HG21	1.84	0.58
25:YA:807:U:O2'	25:YA:2060:A:N1	2.37	0.58
28:YE:78:LEU:HD23	28:YE:79:ARG:HD2	1.86	0.58
29:YF:138:GLU:O	29:YF:141:ALA:HB3	2.03	0.58
32:YI:82:ARG:HD3	32:YI:146:ALA:HB3	1.86	0.58
32:YI:98:ALA:HB2	32:YI:111:PRO:HB3	1.85	0.58
34:YO:64:ARG:HG2	34:YO:79:PHE:CG	2.38	0.58
38:YS:67:ARG:NH1	38:YS:67:ARG:CB	2.64	0.58
45:YZ:52:SER:O	45:YZ:52:SER:OG	2.17	0.58
1:QA:243:A:H4'	1:QA:244:U:O5'	2.03	0.58
8:QH:121:ASP:N	8:QH:121:ASP:OD1	2.35	0.58
13:QM:78:ILE:HG23	13:QM:92:HIS:CD2	2.39	0.58
25:RA:857:C:OP2	46:R0:77:ARG:NH2	2.31	0.58
28:RE:51:PHE:CD2	28:RE:52:LEU:HG	2.39	0.58
33:RN:13:TRP:O	33:RN:135:PRO:HD2	2.03	0.58
2:XB:93:VAL:HG11	2:XB:97:TRP:HD1	1.68	0.58
25:YA:955:C:OP2	36:YQ:14:ARG:HD2	2.03	0.58
25:YA:1012:U:O4	33:YN:25:ARG:HA	2.04	0.58
25:YA:2372:G:O2'	52:Y6:46:HIS:NE2	2.25	0.58
28:YE:111:ARG:NE	28:YE:160:TYR:HE1	2.01	0.58
29:YF:160:ASN:OD1	29:YF:162:LEU:HB2	2.04	0.58
30:YG:113:ARG:HG2	50:Y4:34:GLU:OE2	2.03	0.58
31:YH:4:ILE:HD13	31:YH:4:ILE:H	1.68	0.58
35:YP:39:LYS:HG3	35:YP:45:LEU:HD22	1.86	0.58
40:YU:90:VAL:CG2	41:YV:39:LEU:HB3	2.31	0.58
45:YZ:126:VAL:HG12	45:YZ:163:LEU:HA	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:Y2:15:LYS:H	48:Y2:67:LYS:CE	2.17	0.58
13:QM:58:GLU:O	13:QM:62:ASN:ND2	2.31	0.58
25:RA:1314:C:N4	25:RA:1338:G:H1	2.02	0.58
31:RH:85:LYS:HA	31:RH:86:GLU:OE1	2.04	0.58
33:RN:54:VAL:HB	33:RN:122:VAL:HG22	1.85	0.58
36:RQ:47:ILE:CD1	36:RQ:70:PRO:HD3	2.34	0.58
50:R4:15:ILE:HG22	50:R4:19:GLY:O	2.03	0.58
1:XA:28:G:O2'	1:XA:296:U:OP1	2.21	0.58
1:XA:1320:C:N4	19:XS:36:ARG:HG3	2.19	0.58
25:YA:141:A:C8	25:YA:1408:C:H1'	2.37	0.58
25:YA:573:G:OP2	41:YV:78:LYS:NZ	2.36	0.58
25:YA:1796:U:H2'	25:YA:1797:C:H6	1.68	0.58
25:YA:2599:G:OP2	27:YD:236:GLY:CA	2.52	0.58
27:YD:71:ASP:HB3	27:YD:103:ARG:HH22	1.68	0.58
37:YR:24:GLN:OE1	37:YR:36:THR:HG21	2.04	0.58
37:YR:117:VAL:HG22	37:YR:118:GLU:H	1.68	0.58
48:Y2:21:LEU:O	48:Y2:25:VAL:HG23	2.04	0.58
7:QG:26:PHE:CE2	7:QG:30:ILE:HD11	2.38	0.57
13:QM:120:LYS:O	13:QM:121:LYS:HB2	2.04	0.57
25:RA:27:G:H22	25:RA:512:G:H1'	1.69	0.57
27:RD:24:ILE:HD11	27:RD:91:ARG:HD2	1.85	0.57
28:RE:134:ILE:C	28:RE:134:ILE:HD12	2.28	0.57
30:RG:179:PRO:HB3	50:R4:38:LYS:HD2	1.86	0.57
32:RI:5:LEU:HD13	32:RI:17:GLN:HB3	1.86	0.57
36:RQ:66:ILE:HA	36:RQ:104:PHE:HA	1.85	0.57
50:R4:38:LYS:C	50:R4:40:HIS:H	2.07	0.57
1:XA:455:C:H42	1:XA:477:G:H1	1.52	0.57
1:XA:559:A:OP1	5:XE:126:ARG:NH2	2.37	0.57
1:XA:1122:U:O4	1:XA:1123:A:N6	2.37	0.57
12:XL:83:VAL:HG22	12:XL:84:LEU:H	1.69	0.57
17:XQ:55:ASP:HA	17:XQ:79:SER:HA	1.85	0.57
19:XS:80:TYR:O	19:XS:82:GLY:N	2.36	0.57
25:YA:592:G:N2	54:Y8:4:MET:CE	2.62	0.57
25:YA:2392:A:H2	25:YA:2424:C:H42	1.52	0.57
28:YE:63:LEU:HD13	28:YE:65:GLY:H	1.68	0.57
28:YE:72:VAL:O	28:YE:73:GLU:O	2.21	0.57
39:YT:24:PRO:HA	39:YT:49:VAL:HG13	1.85	0.57
1:QA:1028:C:H42	1:QA:1033:G:H1	1.51	0.57
3:QC:134:ILE:HG23	3:QC:151:VAL:HB	1.85	0.57
25:RA:856:C:O2'	25:RA:857:C:OP1	2.21	0.57
27:RD:35:LYS:HZ1	27:RD:104:TYR:HB2	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:RE:111:ARG:NE	28:RE:160:TYR:HE1	2.01	0.57
31:RH:4:ILE:HD13	31:RH:4:ILE:H	1.68	0.57
31:RH:84:SER:O	31:RH:133:VAL:O	2.22	0.57
44:RY:95:LYS:NZ	44:RY:99:CYS:O	2.37	0.57
50:R4:15:ILE:HG22	50:R4:20:ASN:HA	1.86	0.57
50:R4:27:THR:O	50:R4:28:LYS:HB3	2.03	0.57
51:R5:55:ARG:NH1	51:R5:58:LEU:HD11	2.19	0.57
1:XA:521:G:H4'	12:XL:73:GLU:HG3	1.86	0.57
4:XD:13:ARG:HA	4:XD:33:MET:HE3	1.85	0.57
25:YA:1063:G:H22	25:YA:1076:C:H1'	1.69	0.57
25:YA:1077:A:H5'	25:YA:1078:U:H5"	1.85	0.57
25:YA:1754:C:H5'	39:YT:101:PHE:CE2	2.38	0.57
25:YA:2031:A:N3	25:YA:2455:G:O2'	2.34	0.57
29:YF:53:THR:C	29:YF:55:GLY:H	2.10	0.57
38:YS:95:HIS:CG	38:YS:96:GLY:H	2.21	0.57
25:RA:27:G:H1'	25:RA:513:A:N6	2.20	0.57
25:RA:997:G:OP1	40:RU:93:LYS:HD3	2.05	0.57
30:RG:3:LEU:HD11	50:R4:25:TYR:CZ	2.33	0.57
31:RH:124:GLU:HB3	31:RH:132:ARG:HG3	1.84	0.57
42:RW:71:VAL:HA	42:RW:107:LEU:HD12	1.86	0.57
50:R4:3:GLU:HG3	50:R4:4:GLY:N	2.19	0.57
50:R4:39:CYS:O	50:R4:40:HIS:HB2	2.03	0.57
4:XD:9:CYS:HB3	4:XD:32:ALA:HB2	1.87	0.57
25:YA:1085:A:O2'	25:YA:1086:A:OP1	2.21	0.57
25:YA:1937:A:O2'	25:YA:1939:U:OP2	2.19	0.57
28:YE:63:LEU:HD12	28:YE:65:GLY:H	1.69	0.57
28:YE:116:VAL:CG2	28:YE:122:PHE:CD2	2.86	0.57
28:YE:203:LYS:HE3	28:YE:204:ALA:HB2	1.86	0.57
31:YH:127:GLU:HG2	31:YH:128:PRO:CG	2.32	0.57
35:YP:59:LEU:HA	35:YP:61:ARG:NE	2.20	0.57
42:YW:73:ALA:HB3	42:YW:106:ILE:HD13	1.84	0.57
48:Y2:16:LEU:O	48:Y2:17:SER:HB3	2.04	0.57
1:QA:196:A:OP1	20:QT:68:LYS:NZ	2.37	0.57
25:RA:242:G:H5"	54:R8:62:LEU:CD2	2.34	0.57
25:RA:1543:A:O2'	25:RA:1544:C:H3'	2.03	0.57
32:RI:55:ALA:HA	32:RI:58:LEU:HB3	1.85	0.57
1:XA:1002:G:H2'	1:XA:1003:G:H8	1.69	0.57
1:XA:1392:G:H21	1:XA:1502:A:H8	1.52	0.57
12:XL:33:ARG:O	12:XL:85:ILE:HG22	2.03	0.57
25:YA:222:A:H3'	25:YA:421:U:H5'	1.85	0.57
25:YA:528:A:C2	25:YA:2042:A:H2'	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1567:A:OP2	27:YD:84:TYR:OH	2.15	0.57
25:YA:2286:A:H4'	25:YA:2287:A:O4'	2.05	0.57
27:YD:35:LYS:HG2	27:YD:64:ILE:CA	2.34	0.57
31:YH:3:ARG:HA	31:YH:3:ARG:HE	1.69	0.57
38:YS:26:LEU:CD2	38:YS:87:PHE:HD1	2.17	0.57
54:Y8:46:ARG:O	54:Y8:47:LYS:HB3	2.03	0.57
1:QA:973:G:OP1	10:QJ:57:LYS:NZ	2.37	0.57
13:QM:14:ARG:N	13:QM:44:ARG:HD3	2.18	0.57
28:RE:51:PHE:HD2	28:RE:52:LEU:HG	1.68	0.57
30:RG:22:ARG:HH21	30:RG:171:ALA:HB1	1.69	0.57
30:RG:68:PRO:HB2	30:RG:90:LEU:HD12	1.86	0.57
31:RH:125:VAL:HA	31:RH:126:PRO:CB	2.29	0.57
36:RQ:55:VAL:HG22	36:RQ:56:ARG:N	2.18	0.57
44:RY:76:CYS:SG	44:RY:77:PRO:HD2	2.45	0.57
54:R8:30:ARG:O	54:R8:31:HIS:CB	2.50	0.57
54:R8:53:PRO:CD	54:R8:54:GLU:H	2.15	0.57
1:XA:148:G:H2'	1:XA:149:A:C8	2.38	0.57
1:XA:973:G:OP1	10:XJ:57:LYS:NZ	2.37	0.57
25:YA:654(A):G:H8	25:YA:654(A):G:OP2	1.88	0.57
25:YA:1081:U:H3'	25:YA:1082:U:H4'	1.87	0.57
25:YA:2527:C:H5"	55:Y9:30:PRO:HB2	1.86	0.57
27:YD:263:ARG:HB2	27:YD:263:ARG:HH11	1.68	0.57
28:YE:6:GLY:HA3	28:YE:26:ILE:HD11	1.85	0.57
28:YE:9:VAL:HG11	39:YT:7:ILE:HG22	1.85	0.57
31:YH:2:SER:O	31:YH:3:ARG:C	2.47	0.57
38:YS:5:THR:HG23	38:YS:8:GLU:OE2	2.05	0.57
38:YS:106:ARG:O	38:YS:107:GLU:HB2	2.04	0.57
46:Y0:50:ASN:HB3	46:Y0:63:VAL:HG22	1.86	0.57
1:QA:1014:A:H4'	19:QS:14:HIS:CD2	2.39	0.57
1:QA:1277:C:HO2'	1:QA:1279:A:H8	1.51	0.57
3:QC:134:ILE:HG22	3:QC:168:ALA:HB3	1.87	0.57
7:QG:15:ASP:OD1	7:QG:44:TYR:OH	2.22	0.57
7:QG:79:ARG:HH12	7:QG:82:GLY:HA2	1.69	0.57
27:RD:145:VAL:HG13	27:RD:191:ALA:HB2	1.87	0.57
28:RE:74:PRO:HG2	28:RE:77:ILE:HG23	1.86	0.57
28:RE:102:VAL:HG13	28:RE:172:VAL:CG2	2.34	0.57
31:RH:76:VAL:C	31:RH:78:GLY:H	2.13	0.57
1:XA:1510:U:H2'	1:XA:1511:G:C8	2.39	0.57
3:XC:70:VAL:HG21	3:XC:76:VAL:HG11	1.85	0.57
11:XK:21:ILE:HG13	11:XK:30:VAL:HG12	1.86	0.57
16:XP:20:VAL:HG23	16:XP:35:LYS:HA	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:468:G:N7	53:Y7:39:ARG:NH2	2.52	0.57
25:YA:594:U:P	54:Y8:61:LEU:HD22	2.45	0.57
25:YA:1013:C:H42	25:YA:1149:G:H1	1.52	0.57
25:YA:1805:U:O2	27:YD:50:THR:HB	2.04	0.57
28:YE:41:LYS:HE2	28:YE:41:LYS:HA	1.87	0.57
31:YH:76:VAL:C	31:YH:78:GLY:H	2.13	0.57
37:YR:33:ARG:HH21	51:Y5:55:ARG:HG2	1.69	0.57
1:QA:1321:C:H3'	1:QA:1322:C:H5"	1.85	0.57
12:QL:46:LYS:HG2	12:QL:47:LYS:H	1.70	0.57
25:RA:590:A:OP1	29:RF:95:ARG:NH1	2.37	0.57
28:RE:63:LEU:HD13	28:RE:65:GLY:H	1.68	0.57
28:RE:152:LYS:HB2	33:RN:77:GLY:O	2.05	0.57
30:RG:83:ARG:HG3	30:RG:86:MET:HE3	1.86	0.57
36:RQ:58:PHE:O	36:RQ:59:ARG:C	2.43	0.57
46:R0:18:ALA:O	46:R0:20:ARG:NH1	2.36	0.57
50:R4:37:SER:HB3	50:R4:42:PHE:CE1	2.38	0.57
4:XD:154:ASN:OD1	4:XD:154:ASN:N	2.37	0.57
5:XE:43:LEU:CD2	5:XE:136:MET:HG3	2.35	0.57
25:YA:207:A:H2'	25:YA:208:C:O4'	2.04	0.57
25:YA:1341:U:OP2	25:YA:1394:U:O2'	2.15	0.57
25:YA:2022:U:O2'	25:YA:2617:C:H5'	2.05	0.57
27:YD:28:GLU:O	27:YD:29:PRO:C	2.45	0.57
28:YE:102:VAL:HG13	28:YE:172:VAL:CG2	2.34	0.57
38:YS:72:ALA:O	38:YS:76:LYS:HG3	2.04	0.57
1:QA:359:U:H2'	1:QA:360:A:C8	2.39	0.57
1:QA:895:G:H1	1:QA:904:C:H42	1.51	0.57
6:QF:3:ARG:NH1	6:QF:38:GLU:OE2	2.37	0.57
12:QL:15:ARG:CG	17:QQ:32:TYR:OH	2.53	0.57
13:QM:23:TYR:HB3	13:QM:67:GLU:HG2	1.87	0.57
19:QS:67:VAL:HB	50:R4:59:PHE:CZ	2.40	0.57
27:RD:44:ASN:HB2	27:RD:48:ARG:O	2.05	0.57
53:R7:9:ARG:HH21	53:R7:48:LYS:HB2	1.70	0.57
1:XA:946:A:H2'	1:XA:947:G:C8	2.40	0.57
1:XA:1347:G:H22	1:XA:1374:A:P	2.28	0.57
1:XA:1437:C:N4	1:XA:1464:G:H1	2.03	0.57
25:YA:2275:C:C2	36:YQ:83:MET:HE3	2.39	0.57
29:YF:32:LEU:HD13	29:YF:105:VAL:CG1	2.33	0.57
25:RA:27:G:HO2'	25:RA:28:A:H8	1.50	0.57
25:RA:2250:G:C2	36:RQ:82:ARG:HB3	2.40	0.57
28:RE:41:LYS:HA	28:RE:41:LYS:HE2	1.87	0.57
32:RI:110:ASP:N	32:RI:130:TYR:OH	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:RP:64:LYS:O	35:RP:66:GLY:N	2.38	0.57
37:RR:117:VAL:O	37:RR:118:GLU:HB2	2.04	0.57
41:RV:52:VAL:HG21	41:RV:55:ALA:HB3	1.87	0.57
2:XB:96:ARG:HD3	2:XB:148:TYR:HE1	1.70	0.57
9:XI:15:ALA:HB2	9:XI:65:VAL:HG23	1.86	0.57
19:XS:32:LYS:HA	19:XS:50:ALA:HB3	1.86	0.57
20:XT:53:LEU:O	20:XT:57:ARG:NH1	2.38	0.57
25:YA:2185:C:H2'	25:YA:2186:G:H8	1.69	0.57
25:YA:2250:G:C2	36:YQ:82:ARG:HB3	2.40	0.57
25:YA:2683:C:O2'	28:YE:13:ARG:NH2	2.38	0.57
27:YD:25:THR:HG21	27:YD:82:ILE:H	1.70	0.57
27:YD:36:PRO:HB2	27:YD:61:LEU:HG	1.87	0.57
31:YH:77:LYS:HZ3	31:YH:77:LYS:CB	2.11	0.57
36:YQ:60:ARG:NH1	45:YZ:114:GLY:H	2.01	0.57
38:YS:67:ARG:CB	38:YS:67:ARG:HH11	2.17	0.57
1:QA:1152:A:H5"	10:QJ:13:HIS:HD2	1.69	0.57
2:QB:51:LEU:HD23	2:QB:201:ILE:HD12	1.86	0.57
3:QC:73:PRO:HG3	3:QC:105:GLU:HG3	1.87	0.57
4:QD:154:ASN:N	4:QD:154:ASN:OD1	2.37	0.57
19:QS:41:VAL:HB	19:QS:42:PRO:CA	2.34	0.57
25:RA:592:G:H21	54:R8:4:MET:HE2	1.68	0.57
27:RD:241:PRO:O	27:RD:242:ARG:HB2	2.04	0.57
28:RE:63:LEU:HD12	28:RE:65:GLY:H	1.69	0.57
1:XA:1226:C:OP2	13:XM:103:THR:OG1	2.19	0.57
27:YD:25:THR:HG21	27:YD:81:ALA:HB1	1.86	0.57
27:YD:34:VAL:O	27:YD:34:VAL:CG1	2.51	0.57
29:YF:118:ALA:O	29:YF:121:GLY:N	2.33	0.57
30:YG:81:LYS:O	30:YG:82:LEU:HB2	2.04	0.57
34:YO:85:VAL:HG11	34:YO:114:ILE:HD11	1.87	0.57
20:QT:41:ILE:HG22	20:QT:91:LEU:HD12	1.86	0.56
25:RA:50:U:H3'	25:RA:51:G:H5'	1.86	0.56
25:RA:1899:G:N2	25:RA:1902:C:N4	2.49	0.56
25:RA:2747:G:H21	25:RA:2757:A:H62	1.53	0.56
27:RD:108:PRO:HB3	27:RD:143:HIS:CE1	2.40	0.56
1:XA:57:G:H2'	1:XA:58:C:C6	2.40	0.56
1:XA:438:G:H4'	4:XD:123:HIS:CD2	2.40	0.56
1:XA:636:U:H2'	1:XA:637:G:H8	1.70	0.56
1:XA:953:G:H5'	1:XA:965:A:H61	1.70	0.56
1:XA:1191:A:H5"	3:XC:4:LYS:HZ2	1.70	0.56
25:YA:67:U:C2	25:YA:74:A:H2	2.21	0.56
5:QE:69:VAL:CG1	5:QE:71:LEU:CD2	2.82	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:QL:45:PRO:HD3	12:QL:51:ALA:O	2.04	0.56
25:RA:918:A:N3	26:RB:80:U:O2'	2.35	0.56
27:RD:148:GLU:HB2	27:RD:151:LYS:HD2	1.87	0.56
1:XA:538:G:OP1	12:XL:113:ARG:HD2	2.05	0.56
1:XA:636:U:H2'	1:XA:637:G:C8	2.41	0.56
1:XA:1243:C:H42	1:XA:1294:G:H1	1.51	0.56
1:XA:1443:G:H5"	1:XA:1446:A:H2	1.69	0.56
25:YA:2224:G:OP1	27:YD:268:ARG:HD3	2.04	0.56
27:YD:69:ARG:HD3	27:YD:105:ILE:HD11	1.87	0.56
27:YD:92:ILE:HD12	27:YD:104:TYR:CD2	2.39	0.56
31:YH:84:SER:O	31:YH:133:VAL:O	2.22	0.56
36:YQ:37:LEU:HD21	36:YQ:130:LYS:HE3	1.87	0.56
50:Y4:71:ARG:HB2	50:Y4:71:ARG:HH11	1.68	0.56
1:QA:1065:U:O2'	1:QA:1066:C:OP2	2.19	0.56
1:QA:1443:G:N2	39:RT:119:LYS:HB2	2.18	0.56
5:QE:71:LEU:HD21	5:QE:115:VAL:HG22	1.87	0.56
12:QL:58:VAL:O	12:QL:65:GLU:HA	2.06	0.56
20:QT:33:ILE:HD12	20:QT:63:ILE:CG1	2.36	0.56
25:RA:1089:G:H21	25:RA:1102:C:H42	1.52	0.56
28:RE:37:ARG:NE	28:RE:37:ARG:N	2.53	0.56
28:RE:78:LEU:HD23	28:RE:79:ARG:HD2	1.86	0.56
31:RH:2:SER:O	31:RH:3:ARG:C	2.47	0.56
31:RH:3:ARG:HA	31:RH:3:ARG:HE	1.69	0.56
31:RH:77:LYS:HZ3	31:RH:77:LYS:CB	2.03	0.56
38:RS:106:ARG:HA	38:RS:110:LEU:HD11	1.87	0.56
39:RT:111:ARG:C	39:RT:113:LYS:H	2.12	0.56
1:XA:145:G:H1	1:XA:177:C:H42	1.53	0.56
1:XA:156:G:H1	1:XA:165:C:H42	1.52	0.56
1:XA:735:C:H2'	1:XA:736:C:H6	1.70	0.56
13:XM:120:LYS:C	13:XM:121:LYS:HG2	2.29	0.56
25:YA:26:G:H1'	25:YA:515:A:H61	1.69	0.56
25:YA:993:G:OP1	40:YU:50:ARG:NH2	2.38	0.56
25:YA:1113:U:H2'	25:YA:1114:G:C8	2.40	0.56
25:YA:1359:A:H3'	25:YA:1360:A:C5'	2.34	0.56
25:YA:2277:G:OP2	46:Y0:10:THR:HG21	2.05	0.56
25:YA:2318:G:H22	38:YS:2:ALA:N	2.02	0.56
25:YA:2438:U:O3'	25:YA:2439:A:H3'	2.05	0.56
25:YA:2451:A:N1	56:Z6:101:PPU:HE2	2.21	0.56
25:YA:2585:U:H5	56:Z6:101:PPU:O2'	1.88	0.56
27:YD:2:ALA:O	27:YD:3:VAL:HB	2.06	0.56
27:YD:183:ARG:HD2	27:YD:270:ILE:HG12	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:32:PRO:O	28:YE:34:VAL:HG13	2.06	0.56
32:YI:39:ALA:HB1	32:YI:44:LEU:HD13	1.87	0.56
38:YS:32:LEU:O	38:YS:62:LYS:HE2	2.05	0.56
38:YS:103:GLU:O	38:YS:106:ARG:CG	2.52	0.56
39:YT:60:THR:HG22	39:YT:77:PRO:HA	1.86	0.56
41:YV:59:ALA:HB2	41:YV:96:ILE:HD13	1.88	0.56
41:YV:66:ARG:HH11	41:YV:88:ARG:HD3	1.71	0.56
1:QA:664:G:N2	1:QA:741:G:H1	2.02	0.56
2:QB:82:ARG:HA	2:QB:92:TYR:HE2	1.71	0.56
8:QH:102:ARG:NH1	8:QH:105:ARG:NH2	2.53	0.56
12:QL:18:VAL:O	12:QL:19:ARG:HB2	2.04	0.56
25:RA:884:C:N3	25:RA:892:G:C6	2.73	0.56
25:RA:2811:G:OP1	28:RE:61:ARG:CG	2.54	0.56
29:RF:28:ILE:HG22	29:RF:112:MET:HB3	1.88	0.56
37:RR:67:LEU:HD13	37:RR:76:VAL:HG21	1.86	0.56
44:RY:81:LYS:HZ3	44:RY:98:VAL:HG11	1.69	0.56
50:R4:50:VAL:O	50:R4:51:ASP:C	2.48	0.56
51:R5:55:ARG:HD3	51:R5:56:LYS:N	2.21	0.56
25:YA:898:C:H2'	25:YA:899:A:H5'	1.87	0.56
25:YA:1688:U:O2	25:YA:1700:A:H5"	2.05	0.56
26:YB:90:C:OP1	36:YQ:16:ARG:HG2	2.05	0.56
31:YH:6:ARG:C	31:YH:8:PRO:HD2	2.30	0.56
31:YH:125:VAL:HG12	31:YH:126:PRO:CG	2.34	0.56
36:YQ:79:LEU:O	36:YQ:79:LEU:CG	2.52	0.56
48:Y2:31:GLU:O	48:Y2:35:LEU:HG	2.05	0.56
12:QL:82:VAL:HG23	12:QL:106:ASP:OD2	2.04	0.56
17:QQ:90:ILE:O	17:QQ:94:ASN:ND2	2.38	0.56
25:RA:1543:A:H1'	25:RA:1545:A:O4'	2.05	0.56
28:RE:117:MET:HG3	28:RE:117:MET:O	2.06	0.56
30:RG:6:ALA:HB3	50:R4:23:GLU:HG2	1.88	0.56
36:RQ:79:LEU:O	36:RQ:79:LEU:CG	2.52	0.56
41:RV:44:LYS:HE2	41:RV:45:THR:H	1.70	0.56
50:R4:37:SER:C	50:R4:39:CYS:N	2.62	0.56
54:R8:52:LYS:H	54:R8:53:PRO:HD2	1.66	0.56
16:XP:20:VAL:HG21	16:XP:32:TYR:CD2	2.40	0.56
20:XT:33:ILE:HG23	20:XT:63:ILE:HG12	1.88	0.56
25:YA:575:A:OP2	25:YA:2055:C:N4	2.32	0.56
26:YB:12:C:H2'	46:Y0:73:GLY:HA3	1.87	0.56
26:YB:44:G:H1'	26:YB:47:C:N4	2.20	0.56
28:YE:74:PRO:HG2	28:YE:77:ILE:HG23	1.87	0.56
31:YH:153:LYS:HA	31:YH:153:LYS:HZ3	1.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:YP:65:ARG:O	35:YP:68:GLN:NE2	2.38	0.56
38:YS:14:VAL:HG13	38:YS:15:ARG:N	2.21	0.56
42:YW:71:VAL:HA	42:YW:107:LEU:HD12	1.87	0.56
52:Y6:25:LYS:HE2	52:Y6:27:LYS:HE3	1.87	0.56
1:QA:1306:A:H1'	1:QA:1332:A:C2	2.40	0.56
12:QL:83:VAL:HG22	12:QL:84:LEU:H	1.70	0.56
12:QL:126:LYS:HB2	12:QL:126:LYS:HZ3	1.71	0.56
28:RE:183:LEU:N	28:RE:183:LEU:HD12	2.20	0.56
2:XB:84:GLU:HB3	2:XB:219:VAL:HG21	1.86	0.56
12:XL:111:LYS:O	12:XL:112:ASP:HB2	2.06	0.56
25:YA:263:C:H2'	25:YA:264:C:O4'	2.05	0.56
25:YA:1036:G:OP1	31:YH:59:ARG:HB2	2.05	0.56
28:YE:37:ARG:NE	28:YE:37:ARG:N	2.54	0.56
28:YE:77:ILE:O	28:YE:78:LEU:C	2.48	0.56
48:Y2:41:ILE:HD11	48:Y2:44:LEU:CG	2.36	0.56
10:QJ:5:ARG:HG3	10:QJ:71:LEU:HD11	1.87	0.56
28:RE:32:PRO:O	28:RE:34:VAL:HG13	2.06	0.56
28:RE:64:LYS:C	28:RE:66:HIS:H	2.12	0.56
50:R4:41:PRO:O	50:R4:42:PHE:CB	2.53	0.56
50:R4:48:ARG:O	50:R4:50:VAL:N	2.38	0.56
54:R8:50:LEU:HD12	54:R8:51:ALA:H	1.70	0.56
1:XA:1004:A:H1'	1:XA:1036:G:H1	1.70	0.56
1:XA:1076:C:N4	1:XA:1081:G:H1	2.02	0.56
2:XB:21:ARG:O	2:XB:23:ARG:HD3	2.05	0.56
3:XC:174:PRO:HD2	3:XC:182:ILE:HD11	1.88	0.56
4:XD:11:LEU:HD13	4:XD:66:ARG:HG2	1.87	0.56
7:XG:54:THR:O	7:XG:56:GLN:N	2.39	0.56
27:YD:259:THR:O	27:YD:260:ARG:C	2.49	0.56
28:YE:64:LYS:C	28:YE:66:HIS:H	2.12	0.56
28:YE:183:LEU:HD12	28:YE:183:LEU:N	2.20	0.56
28:YE:195:LEU:HD12	28:YE:196:VAL:H	1.71	0.56
29:YF:197:ASP:O	29:YF:199:TRP:N	2.38	0.56
54:Y8:50:LEU:HD12	54:Y8:51:ALA:H	1.70	0.56
8:QH:106:GLY:O	8:QH:122:ARG:NH2	2.36	0.56
43:RX:31:HIS:CD2	43:RX:32:PRO:HD2	2.41	0.56
1:XA:243:A:H4'	1:XA:244:U:O5'	2.05	0.56
1:XA:501:C:OP1	12:XL:117:ARG:NH2	2.33	0.56
1:XA:560:U:H4'	1:XA:561:U:H5"	1.87	0.56
1:XA:690:G:H2'	1:XA:691:G:O4'	2.05	0.56
1:XA:1446:A:O2'	1:XA:1447:G:O5'	2.23	0.56
25:YA:1782:C:H1'	25:YA:2609:U:H5"	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:YB:50:G:P	38:YS:63:THR:HG23	2.45	0.56
27:YD:239:ARG:O	27:YD:240:ALA:HB2	2.05	0.56
29:YF:155:LEU:CD1	29:YF:174:VAL:HG13	2.32	0.56
31:YH:126:PRO:CG	31:YH:127:GLU:N	2.65	0.56
38:YS:5:THR:OG1	38:YS:7:TYR:HB3	2.06	0.56
28:RE:77:ILE:O	28:RE:78:LEU:C	2.48	0.56
28:RE:203:LYS:HE3	28:RE:204:ALA:HB2	1.86	0.56
31:RH:59:ARG:HH11	31:RH:59:ARG:CG	2.19	0.56
36:RQ:37:LEU:HD21	36:RQ:130:LYS:HE3	1.87	0.56
3:XC:9:GLY:HA2	3:XC:12:LEU:HD23	1.87	0.56
20:XT:29:LYS:HD2	20:XT:66:ALA:HB2	1.87	0.56
25:YA:780:G:N2	25:YA:783:A:H62	2.01	0.56
25:YA:2712:U:HO2'	25:YA:2712(A):A:P	2.28	0.56
27:YD:94:LEU:C	27:YD:94:LEU:HD13	2.30	0.56
28:YE:55:ASN:C	28:YE:57:LYS:N	2.62	0.56
45:YZ:125:LEU:HG	45:YZ:164:ALA:HB3	1.87	0.56
48:Y2:15:LYS:H	48:Y2:67:LYS:NZ	2.02	0.56
1:QA:371:G:H1	1:QA:390:C:H42	1.54	0.56
1:QA:1317:C:N3	19:QS:37:ARG:NH2	2.54	0.56
2:QB:71:VAL:HG12	2:QB:93:VAL:HB	1.88	0.56
22:QV:58:A:O2'	22:QV:60:U:OP2	2.22	0.56
25:RA:2404:C:O3'	35:RP:77:ARG:NH2	2.38	0.56
25:RA:2821:A:OP2	28:RE:110:GLY:CA	2.54	0.56
33:RN:42:TRP:HA	33:RN:48:MET:HE1	1.88	0.56
36:RQ:12:GLN:OE1	36:RQ:72:LYS:HD2	2.06	0.56
50:R4:48:ARG:HH12	50:R4:52:THR:HG22	1.71	0.56
1:XA:1152:A:OP1	10:XJ:68:HIS:NE2	2.39	0.56
1:XA:1162:C:N4	1:XA:1174:G:H1	2.00	0.56
3:XC:130:VAL:O	3:XC:134:ILE:HG12	2.06	0.56
4:XD:122:ARG:NH1	4:XD:134:ASP:O	2.39	0.56
12:XL:79:GLU:O	12:XL:79:GLU:HG2	2.05	0.56
25:YA:1012:U:O2'	25:YA:1013:C:OP2	2.19	0.56
28:YE:117:MET:HG3	28:YE:117:MET:O	2.06	0.56
34:YO:97:ARG:HA	34:YO:117:LEU:HD22	1.88	0.56
35:YP:87:ASP:HB3	35:YP:105:LEU:HD21	1.88	0.56
39:YT:29:ARG:HB2	39:YT:46:GLU:HG3	1.88	0.56
40:YU:92:ARG:HH21	41:YV:10:LYS:HG2	1.69	0.56
44:YY:95:LYS:HB3	44:YY:100:ALA:HA	1.87	0.56
1:QA:757:U:O2'	1:QA:879:C:O2	2.23	0.55
6:QF:69:GLU:CD	6:QF:69:GLU:H	2.15	0.55
9:QI:95:LYS:HZ1	9:QI:96:LEU:HD13	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:QL:111:LYS:O	12:QL:112:ASP:HB2	2.05	0.55
25:RA:1754:C:H5"	39:RT:113:LYS:HE3	1.88	0.55
31:RH:6:ARG:C	31:RH:8:PRO:HD2	2.30	0.55
50:R4:9:LEU:H	50:R4:27:THR:HG22	1.71	0.55
50:R4:61:ARG:C	50:R4:63:TYR:H	2.14	0.55
52:R6:36:LEU:HB2	52:R6:50:ARG:HA	1.89	0.55
1:XA:80:G:O6	1:XA:88:C:N4	2.39	0.55
1:XA:686:U:H1'	11:XK:42:TRP:HE1	1.71	0.55
12:XL:58:VAL:O	12:XL:65:GLU:HA	2.05	0.55
12:XL:83:VAL:HG22	12:XL:84:LEU:N	2.21	0.55
25:YA:581:C:H2'	25:YA:582:G:H8	1.71	0.55
25:YA:1448:G:O2'	25:YA:1528:A:N6	2.40	0.55
25:YA:1818:U:O2'	27:YD:154:LYS:O	2.20	0.55
29:YF:198:ALA:C	29:YF:200:GLU:N	2.62	0.55
29:YF:198:ALA:CA	29:YF:201:VAL:HG12	2.34	0.55
33:YN:56:ASN:N	33:YN:125:GLY:O	2.22	0.55
39:YT:94:ALA:O	39:YT:95:ARG:HB2	2.06	0.55
45:YZ:121:HIS:ND1	45:YZ:123:ASP:O	2.39	0.55
1:QA:1163:C:H42	1:QA:1173:G:H1	1.53	0.55
1:QA:1189:C:OP1	10:QJ:51:ARG:NH2	2.39	0.55
10:QJ:16:LEU:HD23	10:QJ:94:VAL:HG13	1.88	0.55
10:QJ:78:ASN:O	10:QJ:81:THR:OG1	2.24	0.55
25:RA:1111:A:HO2'	31:RH:2:SER:HG	1.51	0.55
25:RA:2419:U:O4	54:R8:30:ARG:CZ	2.55	0.55
26:RB:12:C:H2'	46:R0:73:GLY:HA3	1.89	0.55
34:RO:78:ARG:HH21	39:RT:103:ARG:NH2	2.03	0.55
42:RW:86:LEU:HD22	42:RW:96:ILE:HD11	1.88	0.55
48:R2:46:GLN:O	48:R2:47:ASN:CB	2.53	0.55
1:XA:1333:A:H2'	1:XA:1334:G:O4'	2.07	0.55
5:XE:50:GLU:HB3	5:XE:53:LEU:HD13	1.88	0.55
10:XJ:34:VAL:HG22	10:XJ:74:ILE:HG22	1.89	0.55
25:YA:848:G:H2'	25:YA:849:A:C8	2.41	0.55
25:YA:1142(A):A:H4'	33:YN:25:ARG:HH22	1.71	0.55
25:YA:1287:A:N7	37:YR:107:ASP:HB2	2.20	0.55
25:YA:1799:G:O2'	27:YD:270:ILE:HD11	2.06	0.55
39:YT:3:ARG:HG3	39:YT:7:ILE:HG12	1.88	0.55
1:QA:261:U:OP2	20:QT:79:ARG:NH2	2.40	0.55
20:QT:36:LEU:HD12	20:QT:55:ILE:HG23	1.87	0.55
25:RA:38:A:H2'	25:RA:39:C:C6	2.41	0.55
25:RA:2138:C:H42	25:RA:2153:G:H1	1.53	0.55
28:RE:4:ILE:HD13	28:RE:5:LEU:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:XB:158:LEU:HD13	2:XB:182:ILE:HD11	1.89	0.55
15:XO:18:PHE:CE1	15:XO:21:ASP:HB2	2.41	0.55
31:YH:152:ARG:C	31:YH:153:LYS:HE2	2.32	0.55
36:YQ:12:GLN:OE1	36:YQ:72:LYS:HD2	2.06	0.55
36:YQ:25:ASP:N	36:YQ:102:VAL:HG23	2.22	0.55
38:YS:59:LYS:CG	38:YS:60:GLY:H	2.11	0.55
48:Y2:43:GLN:O	48:Y2:44:LEU:CG	2.54	0.55
1:QA:923:A:N6	1:QA:1392:G:O6	2.40	0.55
2:QB:5:ILE:HD12	2:QB:224:GLN:HG2	1.88	0.55
2:QB:101:MET:HE2	2:QB:108:ILE:HG21	1.88	0.55
30:RG:112:PRO:HB3	50:R4:37:SER:OG	2.06	0.55
31:RH:152:ARG:C	31:RH:153:LYS:HE2	2.32	0.55
34:RO:1:MET:HE3	34:RO:67:LYS:HG2	1.88	0.55
36:RQ:78:PRO:O	46:R0:5:LYS:HE2	2.07	0.55
36:RQ:135:ASP:CG	45:RZ:81:ARG:HH12	2.14	0.55
39:RT:26:ASP:O	39:RT:49:VAL:HG12	2.07	0.55
44:RY:81:LYS:HB2	44:RY:96:ILE:HG22	1.88	0.55
47:R1:53:VAL:HG11	47:R1:90:ILE:HD11	1.88	0.55
54:R8:63:PRO:O	54:R8:64:TYR:HB2	2.07	0.55
1:XA:17:U:H2'	1:XA:18:C:C6	2.42	0.55
1:XA:1054:C:N3	24:XY:34:G:C1'	2.69	0.55
6:XF:97:PHE:HD1	18:XR:31:LEU:HD21	1.70	0.55
25:YA:676:A:N1	25:YA:802:A:N1	2.54	0.55
25:YA:1364:G:C8	47:Y1:2:SER:N	2.74	0.55
25:YA:1674:G:N2	25:YA:1677:A:N1	2.53	0.55
25:YA:1798:U:H5"	27:YD:259:THR:HG22	1.89	0.55
29:YF:29:ASN:N	29:YF:112:MET:HE1	2.19	0.55
36:YQ:64:ILE:HA	36:YQ:106:VAL:CG1	2.33	0.55
39:YT:39:ARG:HG2	39:YT:40:THR:H	1.72	0.55
1:QA:1348:U:N3	1:QA:1374:A:H2	2.05	0.55
8:QH:86:ILE:HG13	8:QH:133:LEU:HD22	1.89	0.55
9:QI:77:ILE:O	9:QI:81:ILE:HG12	2.07	0.55
28:RE:67:PHE:O	28:RE:69:LYS:N	2.39	0.55
32:RI:2:LYS:HA	32:RI:20:ASP:HA	1.87	0.55
44:RY:96:ILE:HG12	44:RY:101:LYS:HB2	1.88	0.55
12:XL:27:LEU:C	12:XL:29:GLY:N	2.64	0.55
25:YA:2757:A:OP1	55:Y9:19:ARG:HA	2.06	0.55
27:YD:35:LYS:CE	27:YD:104:TYR:HB2	2.35	0.55
27:YD:43:ARG:CB	27:YD:54:ARG:HB2	2.37	0.55
27:YD:94:LEU:HD22	27:YD:95:LEU:H	1.69	0.55
35:YP:95:VAL:HG13	35:YP:100:LEU:HD21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:YV:38:LEU:H	41:YV:51:VAL:HG13	1.70	0.55
1:QA:1128:C:H4'	9:QI:16:ARG:HH22	1.72	0.55
22:QV:75:C:OP1	25:RA:2602:A:OP1	2.25	0.55
25:RA:1045:A:O2'	25:RA:1046:A:OP2	2.20	0.55
25:RA:1093:G:P	31:RH:170:ARG:HD2	2.47	0.55
26:RB:113:C:O2'	38:RS:46:VAL:HG13	2.07	0.55
28:RE:3:GLY:HA3	28:RE:81:ILE:HD12	1.88	0.55
32:RI:76:THR:HG21	32:RI:141:LYS:HE3	1.87	0.55
36:RQ:25:ASP:N	36:RQ:102:VAL:HG23	2.21	0.55
43:RX:60:ARG:NH1	53:R7:47:ARG:HH22	2.04	0.55
44:RY:37:VAL:HG21	44:RY:72:VAL:HG21	1.88	0.55
28:YE:3:GLY:HA3	28:YE:81:ILE:HD12	1.88	0.55
28:YE:4:ILE:HD13	28:YE:5:LEU:H	1.71	0.55
31:YH:26:VAL:CG1	31:YH:27:LYS:N	2.63	0.55
32:YI:11:ASN:O	32:YI:12:LEU:HB2	2.07	0.55
38:YS:107:GLU:N	38:YS:110:LEU:HD11	2.22	0.55
52:Y6:28:ARG:HB3	52:Y6:30:THR:H	1.71	0.55
1:QA:565:U:H5"	1:QA:566:G:H2'	1.89	0.55
12:QL:79:GLU:HG2	12:QL:79:GLU:O	2.06	0.55
28:RE:195:LEU:HD12	28:RE:196:VAL:H	1.71	0.55
31:RH:151:ILE:C	31:RH:152:ARG:O	2.49	0.55
41:RV:7:THR:HG23	41:RV:22:VAL:HG11	1.88	0.55
43:RX:25:LYS:HD3	43:RX:80:ILE:HD11	1.89	0.55
1:XA:1000:A:H2'	1:XA:1001:G:C8	2.41	0.55
2:XB:82:ARG:NH1	2:XB:86:GLU:OE2	2.40	0.55
8:XH:86:ILE:HG22	8:XH:93:VAL:HG21	1.89	0.55
25:YA:1061:U:H3'	25:YA:1062:G:H5"	1.88	0.55
28:YE:20:ALA:O	28:YE:21:VAL:CG2	2.48	0.55
29:YF:28:ILE:HD12	29:YF:28:ILE:O	2.06	0.55
32:YI:92:VAL:HG13	32:YI:120:ILE:HG23	1.87	0.55
38:YS:18:ILE:C	38:YS:19:LYS:O	2.44	0.55
46:Y0:27:GLU:HB2	46:Y0:69:PHE:HD1	1.72	0.55
35:RP:106:LEU:O	35:RP:107:LYS:HB2	2.07	0.55
40:RU:52:ARG:HA	40:RU:55:ARG:HG3	1.88	0.55
44:RY:95:LYS:CB	44:RY:100:ALA:HA	2.36	0.55
1:XA:24:U:H2'	1:XA:25:C:H6	1.71	0.55
1:XA:266:G:O2'	1:XA:267:C:OP2	2.18	0.55
11:XK:34:ASP:OD1	11:XK:38:ASN:N	2.39	0.55
28:YE:21:VAL:HG23	28:YE:22:PRO:HD3	1.89	0.55
28:YE:53:PRO:O	28:YE:74:PRO:HA	2.07	0.55
29:YF:24:LEU:HB3	29:YF:115:ALA:HB2	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:YH:143:GLN:HE21	31:YH:143:GLN:C	2.15	0.55
36:YQ:21:THR:O	36:YQ:22:LYS:O	2.25	0.55
39:YT:26:ASP:O	39:YT:49:VAL:HG12	2.07	0.55
48:Y2:31:GLU:HB2	48:Y2:53:LEU:HD11	1.89	0.55
12:QL:83:VAL:HG22	12:QL:84:LEU:N	2.21	0.55
14:QN:48:ALA:HB2	14:QN:53:LEU:HD12	1.89	0.55
25:RA:1050:A:H2'	25:RA:1051:G:O4'	2.06	0.55
25:RA:1490:A:O2'	27:RD:99:ASP:OD1	2.25	0.55
26:RB:37:C:O2	38:RS:95:HIS:NE2	2.39	0.55
27:RD:35:LYS:NZ	27:RD:104:TYR:HB2	2.22	0.55
29:RF:197:ASP:N	29:RF:197:ASP:OD1	2.40	0.55
31:RH:143:GLN:HE21	31:RH:143:GLN:C	2.15	0.55
50:R4:36:CYS:SG	50:R4:39:CYS:CB	2.95	0.55
1:XA:974:A:OP2	14:XN:41:ARG:NH1	2.39	0.55
1:XA:975:A:H5'	1:XA:975:A:H8	1.71	0.55
3:XC:150:LYS:HE2	3:XC:152:ILE:HD11	1.88	0.55
9:XI:16:ARG:HB2	9:XI:64:THR:HB	1.89	0.55
12:XL:28:LYS:O	12:XL:29:GLY:C	2.50	0.55
12:XL:62:SER:C	12:XL:64:TYR:H	2.14	0.55
25:YA:996:A:H4'	40:YU:92:ARG:HE	1.72	0.55
25:YA:1359:A:C3'	25:YA:1360:A:C5'	2.85	0.55
25:YA:1496:A:H8	25:YA:1577:C:O2'	1.89	0.55
25:YA:1952:A:C2	34:YO:22:ILE:HG23	2.41	0.55
28:YE:67:PHE:O	28:YE:69:LYS:N	2.39	0.55
35:YP:65:ARG:NE	54:Y8:15:LYS:HB2	2.22	0.55
41:YV:34:GLU:O	41:YV:36:PRO:HD3	2.06	0.55
42:YW:14:PRO:O	42:YW:17:VAL:N	2.40	0.55
1:QA:996:A:N1	1:QA:1045:C:O2'	2.36	0.55
1:QA:1375:A:H4'	7:QG:29:LYS:HE3	1.88	0.55
9:QI:26:VAL:HG22	9:QI:61:ALA:HB3	1.89	0.55
25:RA:1654:A:C2	28:RE:113:PHE:CD2	2.95	0.55
25:RA:2277:G:OP2	46:R0:10:THR:HG21	2.07	0.55
25:RA:2391:G:OP2	54:R8:32:LEU:HD13	2.06	0.55
25:RA:2414:G:H21	35:RP:67:MET:HE3	1.71	0.55
28:RE:21:VAL:HG23	28:RE:22:PRO:HD3	1.89	0.55
31:RH:8:PRO:O	31:RH:9:ILE:HG23	2.07	0.55
54:R8:56:GLU:O	54:R8:57:ARG:C	2.50	0.55
25:YA:1820:U:O2	27:YD:202:LYS:HB3	2.07	0.55
25:YA:2393:A:H5'	35:YP:62:LEU:HB3	1.89	0.55
39:YT:62:THR:HG22	39:YT:75:ILE:HG12	1.89	0.55
41:YV:61:VAL:HG23	41:YV:63:GLY:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:YZ:72:ARG:NH2	45:YZ:97:GLU:O	2.31	0.55
1:QA:411:A:N6	1:QA:413:G:H21	2.05	0.54
10:QJ:13:HIS:CE1	10:QJ:14:LYS:HE3	2.42	0.54
13:QM:62:ASN:OD1	50:R4:49:PHE:HD2	1.89	0.54
22:QV:3:G:HO2'	22:QV:4:G:H8	1.53	0.54
25:RA:2413:G:H21	35:RP:70:GLN:HE22	1.55	0.54
25:RA:2645:G:H3'	25:RA:2646:C:H5'	1.88	0.54
48:R2:48:HIS:C	48:R2:48:HIS:CD2	2.85	0.54
1:XA:1007:C:H2'	1:XA:1008:C:H5"	1.89	0.54
4:XD:92:VAL:O	4:XD:96:LEU:HD22	2.07	0.54
8:XH:54:ASP:N	8:XH:54:ASP:OD1	2.39	0.54
25:YA:229:A:OP1	25:YA:229:A:H4'	2.06	0.54
25:YA:247:G:H4'	25:YA:386:G:C5	2.42	0.54
28:YE:54:GLN:NE2	28:YE:54:GLN:N	2.56	0.54
29:YF:32:LEU:HD12	29:YF:36:VAL:HG23	1.89	0.54
38:YS:36:TYR:HD2	38:YS:52:SER:CB	2.18	0.54
54:Y8:30:ARG:O	54:Y8:31:HIS:CB	2.55	0.54
1:QA:1172:C:H2'	1:QA:1173:G:C8	2.43	0.54
25:RA:389:G:H1	35:RP:70:GLN:HB3	1.72	0.54
25:RA:873:G:N2	25:RA:904:C:N3	2.48	0.54
25:RA:2572:A:C4	28:RE:144:ARG:NH2	2.75	0.54
28:RE:152:LYS:HG2	33:RN:78:TYR:CE1	2.42	0.54
31:RH:12:PRO:O	31:RH:13:LYS:HB2	2.07	0.54
33:RN:30:ILE:HG22	33:RN:34:LEU:HD22	1.88	0.54
33:RN:40:PRO:HB3	40:RU:68:ALA:HB2	1.89	0.54
39:RT:1:MET:O	39:RT:3:ARG:N	2.40	0.54
1:XA:690:G:H22	11:XK:55:LYS:HZ1	1.53	0.54
1:XA:982:U:H5"	14:XN:6:LEU:HD21	1.88	0.54
5:XE:11:ILE:CG2	5:XE:108:ALA:HB2	2.38	0.54
13:XM:7:VAL:HG21	30:YG:113:ARG:O	2.06	0.54
25:YA:581:C:H2'	25:YA:582:G:C8	2.42	0.54
25:YA:1794:U:H2'	25:YA:1795:C:C6	2.41	0.54
25:YA:1996:C:OP1	34:YO:31:LYS:NZ	2.40	0.54
25:YA:2115:G:N2	25:YA:2165:G:N7	2.46	0.54
27:YD:31:LYS:O	27:YD:35:LYS:O	2.24	0.54
27:YD:65:ILE:HD13	27:YD:65:ILE:C	2.33	0.54
28:YE:176:ILE:HG22	28:YE:179:GLU:H	1.72	0.54
30:YG:179:PRO:HG3	50:Y4:38:LYS:HZ1	1.70	0.54
33:YN:39:ARG:HD3	33:YN:48:MET:HE2	1.87	0.54
37:YR:55:ALA:HB2	37:YR:79:LEU:HD13	1.89	0.54
1:QA:382:A:H2'	1:QA:383:A:H8	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:411:A:C4	1:QA:413:G:H1'	2.42	0.54
4:QD:31:CYS:SG	4:QD:33:MET:HB2	2.47	0.54
12:QL:28:LYS:O	12:QL:29:GLY:C	2.50	0.54
25:RA:27:G:N2	25:RA:512:G:H1'	2.22	0.54
25:RA:65:C:O2'	25:RA:456:C:N3	2.40	0.54
25:RA:1086:A:H4'	25:RA:1103:A:N6	2.22	0.54
28:RE:61:ARG:O	28:RE:62:PRO:C	2.50	0.54
31:RH:128:PRO:CD	31:RH:129:THR:N	2.71	0.54
1:XA:67:C:H2'	1:XA:68:G:C8	2.42	0.54
1:XA:129(A):G:O2'	1:XA:189:U:H3'	2.08	0.54
1:XA:1321:C:H5'	13:XM:87:TYR:CE1	2.42	0.54
13:XM:20:THR:C	13:XM:22:ILE:H	2.15	0.54
13:XM:22:ILE:HD12	13:XM:25:ILE:HD12	1.89	0.54
25:YA:530:G:N1	25:YA:2022:U:OP1	2.40	0.54
25:YA:1057:A:H62	25:YA:1086:A:H2'	1.71	0.54
25:YA:2506:U:H1'	56:Z6:101:PPU:HN'3	1.71	0.54
28:YE:15:PHE:CD1	39:YT:81:PRO:CD	2.90	0.54
31:YH:8:PRO:O	31:YH:9:ILE:HG23	2.08	0.54
37:YR:56:LYS:NZ	37:YR:87:TYR:O	2.40	0.54
51:Y5:55:ARG:HG3	51:Y5:57:VAL:H	1.73	0.54
54:Y8:56:GLU:O	54:Y8:57:ARG:C	2.50	0.54
10:QJ:8:LEU:HB3	10:QJ:16:LEU:HD21	1.88	0.54
13:QM:22:ILE:HB	13:QM:25:ILE:HD12	1.89	0.54
25:RA:2115:G:N2	25:RA:2165:G:N7	2.56	0.54
25:RA:2350:C:H5	54:R8:42:ARG:NH1	2.05	0.54
26:RB:33:G:H5'	30:RG:2:PRO:HG3	1.89	0.54
26:RB:52:A:N6	38:RS:33:LYS:HG3	2.22	0.54
31:RH:86:GLU:HG3	31:RH:165:ALA:CB	2.38	0.54
36:RQ:39:PRO:HB3	36:RQ:99:PRO:HD3	1.90	0.54
40:RU:112:ARG:NH2	41:RV:47:VAL:HG13	2.23	0.54
50:R4:65:ASP:O	50:R4:66:SER:HB3	2.07	0.54
1:XA:266:G:C8	1:XA:266:G:C5'	2.91	0.54
1:XA:1502:A:H2	1:XA:1505:G:N1	2.06	0.54
5:XE:12:LEU:HD21	5:XE:14:ARG:HD3	1.89	0.54
10:XJ:33:GLN:HB2	10:XJ:75:ILE:HD11	1.89	0.54
18:XR:31:LEU:HD23	18:XR:31:LEU:H	1.73	0.54
25:YA:27:G:O2'	25:YA:28:A:H8	1.91	0.54
27:YD:155:LEU:HD23	27:YD:177:LEU:CD2	2.36	0.54
28:YE:14:ILE:HG23	28:YE:15:PHE:N	2.22	0.54
31:YH:86:GLU:HG3	31:YH:165:ALA:CB	2.38	0.54
40:YU:76:TYR:CZ	40:YU:80:ILE:HG13	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:Y4:54:GLY:O	50:Y4:59:PHE:HB2	2.07	0.54
1:QA:660:G:H1	1:QA:745:C:H42	1.54	0.54
1:QA:1151:A:H2'	1:QA:1152:A:C8	2.43	0.54
25:RA:994:C:OP2	40:RU:54:LYS:NZ	2.41	0.54
25:RA:1441:G:H2'	25:RA:1442:G:H8	1.72	0.54
25:RA:2619:C:O2'	28:RE:156:MET:CE	2.55	0.54
28:RE:101:ARG:HB3	28:RE:201:THR:OG1	2.08	0.54
36:RQ:21:THR:O	36:RQ:22:LYS:O	2.25	0.54
37:RR:103:ARG:NH1	37:RR:108:GLY:O	2.41	0.54
43:RX:83:VAL:CG1	43:RX:87:GLN:HB2	2.38	0.54
47:R1:58:ILE:HD11	47:R1:86:SER:HB2	1.88	0.54
50:R4:37:SER:HB3	50:R4:42:PHE:CD1	2.43	0.54
51:R5:20:ARG:C	51:R5:22:HIS:H	2.14	0.54
9:XI:40:LEU:C	9:XI:42:ARG:H	2.15	0.54
10:XJ:50:ILE:CD1	10:XJ:57:LYS:CG	2.77	0.54
25:YA:551:G:H5'	25:YA:1220:A:H1'	1.90	0.54
27:YD:183:ARG:HG2	27:YD:183:ARG:NH1	2.12	0.54
29:YF:62:ARG:HB3	29:YF:62:ARG:NH1	2.22	0.54
29:YF:129:PHE:O	29:YF:130:ALA:CB	2.55	0.54
31:YH:153:LYS:CE	31:YH:153:LYS:HA	2.38	0.54
46:Y0:20:ARG:O	46:Y0:24:LYS:NZ	2.39	0.54
54:Y8:32:LEU:O	54:Y8:36:LYS:HE3	2.07	0.54
9:QI:121:ARG:NH1	9:QI:122:ALA:O	2.40	0.54
25:RA:102:G:H4'	25:RA:103:A:O5'	2.07	0.54
25:RA:252:G:OP2	35:RP:50:ARG:NH1	2.40	0.54
25:RA:635:C:O2'	25:RA:639:U:OP1	2.26	0.54
28:RE:186:GLY:O	28:RE:188:VAL:N	2.41	0.54
29:RF:181:LEU:HD13	29:RF:186:ILE:HD11	1.89	0.54
45:RZ:108:PRO:HA	45:RZ:142:SER:HA	1.90	0.54
48:R2:42:GLY:O	48:R2:44:LEU:N	2.35	0.54
54:R8:32:LEU:O	54:R8:36:LYS:HE3	2.07	0.54
1:XA:892:A:H2'	1:XA:893:C:C6	2.42	0.54
1:XA:1118:C:P	9:XI:104:ARG:HH11	2.30	0.54
12:XL:6:THR:OG1	12:XL:9:GLN:HG3	2.08	0.54
20:XT:43:LEU:HD13	20:XT:51:GLU:HG3	1.89	0.54
25:YA:1639:U:H2'	25:YA:1640:C:H5"	1.89	0.54
25:YA:2864:G:OP1	39:YT:119:LYS:HD2	2.06	0.54
27:YD:72:LYS:O	27:YD:73:VAL:C	2.51	0.54
27:YD:118:VAL:HG22	27:YD:119:ALA:H	1.72	0.54
27:YD:124:PRO:HB2	27:YD:126:GLN:NE2	2.22	0.54
27:YD:158:ALA:HB3	27:YD:161:THR:HG21	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:206:LEU:O	27:YD:211:ARG:NH1	2.38	0.54
28:YE:14:ILE:HD11	39:YT:14:TYR:CZ	2.41	0.54
33:YN:35:ARG:HB2	33:YN:42:TRP:CH2	2.42	0.54
36:YQ:60:ARG:NH1	45:YZ:114:GLY:N	2.55	0.54
1:QA:607:A:H2'	1:QA:608:A:O4'	2.07	0.54
1:QA:1109:C:H2'	1:QA:1110:A:O4'	2.08	0.54
12:QL:6:THR:O	12:QL:7:ILE:C	2.51	0.54
12:QL:62:SER:C	12:QL:64:TYR:H	2.14	0.54
25:RA:774:A:H2	25:RA:787:U:HO2'	1.52	0.54
28:RE:15:PHE:CE1	39:RT:81:PRO:HD3	2.42	0.54
28:RE:176:ILE:HG22	28:RE:179:GLU:H	1.71	0.54
30:RG:16:ARG:NH2	30:RG:28:VAL:O	2.41	0.54
31:RH:86:GLU:H	31:RH:86:GLU:CD	2.16	0.54
36:RQ:60:ARG:HH12	36:RQ:113:GLN:HE22	1.55	0.54
39:RT:111:ARG:O	39:RT:112:ARG:HG3	2.08	0.54
51:R5:55:ARG:HD3	51:R5:56:LYS:H	1.73	0.54
1:XA:265:G:N2	1:XA:267:C:H5'	2.22	0.54
3:XC:150:LYS:HB3	3:XC:201:TYR:HB2	1.90	0.54
7:XG:20:ASP:HB3	7:XG:23:VAL:HG23	1.89	0.54
7:XG:49:ILE:O	7:XG:53:LYS:HB3	2.08	0.54
16:XP:43:LYS:HG2	16:XP:48:TRP:CE3	2.42	0.54
25:YA:857:C:OP2	46:Y0:77:ARG:NH2	2.40	0.54
26:YB:90:C:OP1	36:YQ:16:ARG:CG	2.56	0.54
27:YD:211:ARG:HD2	27:YD:214:TRP:CZ3	2.43	0.54
29:YF:127:GLU:O	29:YF:129:PHE:N	2.39	0.54
31:YH:91:GLY:O	31:YH:94:TYR:HB2	2.08	0.54
32:YI:21:VAL:HG21	32:YI:25:TYR:HD2	1.73	0.54
1:QA:452:A:O2'	1:QA:453:A:O4'	2.24	0.54
1:QA:1151:A:H2'	1:QA:1152:A:H8	1.73	0.54
1:QA:1346:A:H5"	9:QI:120:ARG:NH1	2.19	0.54
4:QD:150:GLU:HA	4:QD:153:ARG:HG2	1.90	0.54
25:RA:247:G:H4'	25:RA:386:G:C5	2.42	0.54
25:RA:880:G:O6	25:RA:897:C:N3	2.41	0.54
25:RA:943:U:OP2	35:RP:36:LYS:HG2	2.07	0.54
25:RA:1210:A:H5"	25:RA:1210:A:H8	1.72	0.54
31:RH:26:VAL:CG1	31:RH:27:LYS:N	2.64	0.54
32:RI:11:ASN:O	32:RI:12:LEU:HB2	2.08	0.54
36:RQ:134:ARG:HH12	45:RZ:119:GLU:HG3	1.73	0.54
50:R4:51:ASP:O	50:R4:51:ASP:OD1	2.25	0.54
52:R6:15:GLU:CD	52:R6:41:PRO:HB3	2.32	0.54
1:XA:272:C:H2'	1:XA:273:A:C8	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:XL:83:VAL:CG2	12:XL:100:ILE:HG23	2.38	0.54
12:XL:126:LYS:HB2	12:XL:126:LYS:HZ2	1.72	0.54
13:XM:3:ARG:HA	13:XM:9:ILE:CG2	2.37	0.54
25:YA:2633:G:H1'	28:YE:62:PRO:HG2	1.90	0.54
26:YB:8:U:O2'	38:YS:40:ILE:HD13	2.08	0.54
27:YD:130:ALA:C	27:YD:131:LEU:HD12	2.33	0.54
28:YE:51:PHE:O	28:YE:74:PRO:HB3	2.08	0.54
38:YS:13:ARG:HD2	38:YS:13:ARG:O	2.06	0.54
1:QA:662:G:O2'	1:QA:836:G:OP1	2.26	0.54
2:QB:84:GLU:HB3	2:QB:219:VAL:HG21	1.89	0.54
5:QE:7:GLU:N	5:QE:35:GLY:O	2.36	0.54
12:QL:83:VAL:CG2	12:QL:100:ILE:HG23	2.38	0.54
19:QS:5:LEU:HD13	50:R4:67:TYR:HE2	1.71	0.54
25:RA:884:C:C2	25:RA:892:G:N1	2.76	0.54
25:RA:1262:A:C2	51:R5:10:LYS:HE3	2.42	0.54
25:RA:1337:G:OP2	43:RX:73:ARG:NH2	2.40	0.54
25:RA:1952:A:C2	34:RO:22:ILE:HG23	2.43	0.54
27:RD:228:PRO:HD3	27:RD:234:GLY:C	2.33	0.54
28:RE:14:ILE:CG1	28:RE:15:PHE:H	2.08	0.54
28:RE:53:PRO:O	28:RE:74:PRO:HA	2.07	0.54
29:RF:24:LEU:HD23	29:RF:115:ALA:HA	1.89	0.54
31:RH:91:GLY:O	31:RH:94:TYR:HB2	2.08	0.54
39:RT:37:GLY:O	39:RT:39:ARG:N	2.34	0.54
1:XA:686:U:O4	1:XA:703:G:H1'	2.08	0.54
1:XA:1241:G:H2'	1:XA:1242:C:C6	2.42	0.54
3:XC:15:THR:HG23	3:XC:181:ASN:HD22	1.73	0.54
6:XF:50:TYR:OH	18:XR:74:ARG:O	2.20	0.54
10:XJ:4:ILE:HG12	10:XJ:100:THR:HG22	1.89	0.54
10:XJ:9:ARG:HB2	10:XJ:95:GLU:HB3	1.88	0.54
25:YA:242:G:O5'	54:Y8:3:LYS:HE3	2.07	0.54
25:YA:1803:A:C4'	27:YD:259:THR:HG21	2.35	0.54
25:YA:2250:G:C4	36:YQ:82:ARG:HG3	2.43	0.54
27:YD:25:THR:O	27:YD:25:THR:HG23	2.07	0.54
27:YD:80:ALA:HB3	27:YD:94:LEU:HD13	1.88	0.54
30:YG:15:VAL:HG21	30:YG:176:LEU:HD23	1.90	0.54
31:YH:126:PRO:HD2	31:YH:127:GLU:H	1.72	0.54
45:YZ:182:LYS:HG3	45:YZ:183:LEU:CA	2.15	0.54
1:QA:1405:G:OP2	58:QA:1696:PAR:O34	2.25	0.54
25:RA:859:G:N2	25:RA:917:A:OP2	2.41	0.54
25:RA:2749:A:OP1	31:RH:4:ILE:HG22	2.07	0.54
36:RQ:83:MET:H	46:R0:7:LEU:HD12	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:RR:45:ARG:HA	37:RR:95:THR:HG21	1.90	0.54
10:XJ:32:ALA:H	10:XJ:78:ASN:HD21	1.55	0.54
12:XL:6:THR:O	12:XL:7:ILE:C	2.51	0.54
25:YA:286:C:H2'	25:YA:287:C:H6	1.73	0.54
25:YA:2585:U:C5	56:Z6:101:PPU:O2'	2.61	0.54
27:YD:85:ASP:OD2	27:YD:88:ARG:HG2	2.07	0.54
28:YE:101:ARG:HB3	28:YE:201:THR:OG1	2.08	0.54
28:YE:186:GLY:O	28:YE:188:VAL:N	2.40	0.54
31:YH:86:GLU:H	31:YH:86:GLU:CD	2.16	0.54
33:YN:6:PRO:HG3	33:YN:41:ASP:HB2	1.89	0.54
35:YP:49:ARG:HG3	54:Y8:59:LYS:HG2	1.89	0.54
36:YQ:39:PRO:HB3	36:YQ:99:PRO:HD3	1.90	0.54
1:QA:407:G:O4'	4:QD:119:GLN:NE2	2.41	0.53
2:QB:235:SER:O	2:QB:237:ALA:N	2.41	0.53
5:QE:33:VAL:HG12	5:QE:112:LEU:HD12	1.90	0.53
12:QL:11:VAL:HG13	17:QQ:29:HIS:HD2	1.73	0.53
12:QL:119:LYS:C	12:QL:120:TYR:HD1	2.16	0.53
19:QS:10:PHE:HE1	19:QS:16:LEU:HD22	1.73	0.53
25:RA:2116:G:H1	25:RA:2162:G:P	2.29	0.53
28:RE:51:PHE:O	28:RE:74:PRO:HB3	2.08	0.53
28:RE:54:GLN:NE2	28:RE:54:GLN:N	2.55	0.53
31:RH:139:GLN:O	31:RH:143:GLN:HB2	2.09	0.53
1:XA:1086:U:H3	1:XA:1099:G:H22	1.55	0.53
2:XB:82:ARG:HA	2:XB:92:TYR:CE2	2.43	0.53
12:XL:119:LYS:C	12:XL:120:TYR:HD1	2.16	0.53
25:YA:1022:G:N2	25:YA:1023:U:O4	2.38	0.53
25:YA:2123:G:H2'	25:YA:2124:G:C8	2.43	0.53
26:YB:75:G:C5'	45:YZ:36:LYS:HE2	2.37	0.53
27:YD:198:ASN:C	27:YD:198:ASN:HD22	2.16	0.53
31:YH:59:ARG:HH11	31:YH:59:ARG:CG	2.19	0.53
31:YH:125:VAL:HA	31:YH:126:PRO:CB	2.29	0.53
31:YH:151:ILE:C	31:YH:152:ARG:O	2.49	0.53
33:YN:30:ILE:HG23	33:YN:52:VAL:HG11	1.90	0.53
1:QA:1053:G:H5"	1:QA:1055:A:OP1	2.09	0.53
1:QA:1350:A:OP2	9:QI:118:LYS:NZ	2.40	0.53
12:QL:6:THR:OG1	12:QL:9:GLN:HG3	2.08	0.53
15:QO:6:GLU:OE2	15:QO:6:GLU:N	2.35	0.53
25:RA:2133:G:H1'	25:RA:2158:A:H61	1.72	0.53
25:RA:2585:U:H5	56:Z5:101:PPU:HO2'	1.57	0.53
39:RT:33:LYS:HD2	39:RT:82:LEU:HA	1.89	0.53
51:R5:60:VAL:OXT	51:R5:60:VAL:CG1	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:XB:162:ILE:O	2:XB:185:ILE:HG12	2.08	0.53
3:XC:14:ILE:O	3:XC:16:ARG:N	2.35	0.53
25:YA:49:A:C8	25:YA:120:U:C5	2.96	0.53
25:YA:1535:U:H5"	25:YA:1537:C:C4	2.43	0.53
25:YA:2277:G:OP1	36:YQ:85:LYS:HB2	2.08	0.53
25:YA:2477:C:H3'	55:Y9:1:MET:HE3	1.90	0.53
29:YF:147:GLY:O	29:YF:148:LEU:HD23	2.08	0.53
29:YF:197:ASP:O	29:YF:198:ALA:HB3	2.06	0.53
38:YS:25:ARG:HH12	38:YS:42:ASP:CG	2.16	0.53
41:YV:66:ARG:NH1	41:YV:88:ARG:HD3	2.23	0.53
48:Y2:41:ILE:HD11	48:Y2:44:LEU:HB2	1.90	0.53
54:Y8:63:PRO:O	54:Y8:64:TYR:HB2	2.07	0.53
1:QA:1305:G:H5'	21:QU:4:GLY:HA3	1.90	0.53
3:QC:35:GLU:HG2	3:QC:59:ARG:NH2	2.23	0.53
3:QC:84:ILE:HD11	3:QC:88:ARG:HH21	1.73	0.53
10:QJ:22:LYS:HZ1	10:QJ:23:ILE:HG12	1.73	0.53
12:QL:42:THR:HA	12:QL:53:ARG:O	2.08	0.53
25:RA:620:G:H4'	25:RA:621:A:H5"	1.89	0.53
27:RD:44:ASN:HD22	27:RD:44:ASN:N	2.06	0.53
30:RG:65:GLY:O	50:R4:7:PRO:CG	2.56	0.53
31:RH:153:LYS:HA	31:RH:153:LYS:CE	2.38	0.53
34:RO:2:ILE:HD13	34:RO:8:LEU:HD11	1.90	0.53
36:RQ:81:VAL:HG23	46:R0:7:LEU:HD11	1.90	0.53
51:R5:16:ARG:NH1	51:R5:17:ASP:OD1	2.41	0.53
52:R6:25:LYS:CE	54:R8:34:TRP:HZ2	2.20	0.53
54:R8:29:LYS:HB2	54:R8:44:LYS:HG2	1.90	0.53
1:XA:918:A:H2'	1:XA:919:A:C8	2.43	0.53
1:XA:1256:A:OP2	3:XC:26:LYS:NZ	2.39	0.53
1:XA:1392:G:H2'	1:XA:1393:U:H6	1.72	0.53
5:XE:11:ILE:HG21	5:XE:108:ALA:CB	2.38	0.53
12:XL:42:THR:HA	12:XL:53:ARG:O	2.08	0.53
25:YA:860:U:OP2	25:YA:916:G:N1	2.39	0.53
25:YA:1007:C:H4'	33:YN:108:PRO:HD3	1.91	0.53
25:YA:2443:C:OP1	29:YF:68:LYS:CD	2.56	0.53
29:YF:116:ASP:OD2	35:YP:1:MET:N	2.34	0.53
35:YP:61:ARG:HD2	54:Y8:13:ARG:HD2	1.89	0.53
36:YQ:112:GLU:H	36:YQ:112:GLU:CD	2.17	0.53
1:QA:1262:C:H42	1:QA:1273:G:H1	1.55	0.53
4:QD:13:ARG:HA	4:QD:33:MET:HE3	1.90	0.53
5:QE:89:ILE:HG13	5:QE:121:LYS:O	2.09	0.53
13:QM:65:LYS:HE2	50:R4:50:VAL:HG11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:QS:68:GLY:CA	50:R4:68:ARG:CB	2.85	0.53
25:RA:458:G:O2'	25:RA:469:G:O6	2.23	0.53
25:RA:2022:U:O2'	25:RA:2617:C:H5'	2.07	0.53
32:RI:88:ILE:O	32:RI:121:LYS:NZ	2.39	0.53
36:RQ:119:ARG:HG2	36:RQ:119:ARG:NH1	2.19	0.53
41:RV:99:ILE:O	41:RV:101:GLY:N	2.42	0.53
54:R8:52:LYS:N	54:R8:53:PRO:HD2	2.22	0.53
1:XA:855:G:OP2	1:XA:871:U:N3	2.32	0.53
1:XA:1376:U:H2'	1:XA:1377:A:C8	2.44	0.53
25:YA:878:A:N6	25:YA:899:A:O2'	2.40	0.53
25:YA:1441:G:H2'	25:YA:1442:G:H8	1.73	0.53
25:YA:1858:G:O2'	25:YA:1884:A:N6	2.40	0.53
25:YA:2123:G:H1	25:YA:2175:C:N4	2.07	0.53
25:YA:2443:C:OP1	29:YF:68:LYS:HD2	2.09	0.53
25:YA:2777:G:OP2	25:YA:2781:A:O2'	2.24	0.53
28:YE:61:ARG:O	28:YE:62:PRO:C	2.51	0.53
38:YS:74:ALA:HB1	38:YS:107:GLU:HB3	1.89	0.53
45:YZ:124:ILE:HG22	45:YZ:126:VAL:HG13	1.90	0.53
55:Y9:27:CYS:SG	55:Y9:28:GLU:N	2.82	0.53
1:QA:1241:G:H2'	1:QA:1242:C:C6	2.44	0.53
1:QA:1306:A:N6	1:QA:1331:G:H1'	2.22	0.53
5:QE:75:THR:OG1	5:QE:117:ASP:O	2.26	0.53
25:RA:2712:U:O2'	25:RA:2712(A):A:H8	1.91	0.53
29:RF:66:PRO:O	29:RF:68:LYS:N	2.41	0.53
38:RS:106:ARG:HA	38:RS:110:LEU:HD21	1.91	0.53
51:R5:44:THR:O	51:R5:46:CYS:N	2.41	0.53
1:XA:143:A:H2	1:XA:220:G:H1	1.57	0.53
2:XB:60:ASP:O	2:XB:64:ARG:HG2	2.09	0.53
9:XI:111:ARG:NE	9:XI:112:LYS:O	2.38	0.53
12:XL:32:PHE:HE1	12:XL:86:ARG:HG3	1.73	0.53
17:XQ:4:LYS:HE3	17:XQ:6:LEU:HD21	1.90	0.53
19:XS:36:ARG:NH1	19:XS:52:TYR:O	2.42	0.53
25:YA:7:G:H1	25:YA:2896:C:N4	2.04	0.53
25:YA:825:C:O2	35:YP:55:ARG:NH2	2.41	0.53
25:YA:1035:U:H2'	25:YA:1036:G:C8	2.44	0.53
28:YE:47:VAL:O	28:YE:48:GLN:C	2.52	0.53
31:YH:12:PRO:O	31:YH:13:LYS:HB2	2.07	0.53
33:YN:40:PRO:O	40:YU:64:ARG:HD2	2.09	0.53
36:YQ:64:ILE:HG13	45:YZ:178:GLU:OE1	2.09	0.53
39:YT:51:ARG:CG	39:YT:98:LYS:HG3	2.38	0.53
39:YT:112:ARG:O	39:YT:112:ARG:NE	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:Y8:58:ILE:O	54:Y8:61:LEU:HG	2.08	0.53
1:QA:323:U:H5'	20:QT:23:ARG:HB2	1.89	0.53
1:QA:580:U:H2'	1:QA:581:G:O4'	2.09	0.53
1:QA:957:U:H4'	19:QS:79:THR:HB	1.91	0.53
8:QH:51:VAL:HG11	8:QH:60:ARG:HG3	1.90	0.53
9:QI:71:SER:HA	9:QI:74:ILE:HD12	1.90	0.53
19:QS:15:LEU:HD12	19:QS:44:MET:HE1	1.90	0.53
20:QT:26:ASN:CB	20:QT:71:THR:HG23	2.38	0.53
25:RA:910:A:H62	36:RQ:12:GLN:HA	1.72	0.53
25:RA:2123:G:H2'	25:RA:2124:G:C8	2.41	0.53
35:RP:122:PRO:HA	35:RP:141:ALA:HB1	1.90	0.53
36:RQ:112:GLU:CD	36:RQ:112:GLU:H	2.17	0.53
45:RZ:117:LEU:HA	45:RZ:174:VAL:HA	1.91	0.53
48:R2:10:LEU:O	48:R2:13:ALA:N	2.40	0.53
54:R8:58:ILE:O	54:R8:61:LEU:HG	2.08	0.53
1:XA:481:G:O2'	1:XA:482:A:O5'	2.27	0.53
5:XE:41:VAL:HG11	5:XE:113:ALA:HB2	1.85	0.53
25:YA:443:A:H5"	25:YA:444:C:OP1	2.09	0.53
25:YA:1593:G:H2'	25:YA:1594:G:C8	2.44	0.53
27:YD:25:THR:CG2	27:YD:81:ALA:HB1	2.38	0.53
27:YD:35:LYS:CG	27:YD:64:ILE:H	2.15	0.53
31:YH:128:PRO:CD	31:YH:129:THR:N	2.71	0.53
35:YP:49:ARG:HG3	54:Y8:59:LYS:CG	2.39	0.53
25:RA:1427:A:H4'	25:RA:1428:C:O5'	2.07	0.53
25:RA:2808:U:H3	25:RA:2892:A:N6	1.99	0.53
28:RE:1:MET:HE2	28:RE:1:MET:H1	1.73	0.53
32:RI:52:ARG:CA	32:RI:56:LYS:HB3	2.38	0.53
35:RP:65:ARG:NH2	54:R8:46:ARG:HH12	2.07	0.53
44:RY:98:VAL:HG13	44:RY:99:CYS:SG	2.48	0.53
45:RZ:27:VAL:HG13	45:RZ:87:ASP:HB3	1.91	0.53
50:R4:42:PHE:O	50:R4:43:TYR:C	2.51	0.53
1:XA:34:C:H2'	1:XA:35:G:C8	2.43	0.53
1:XA:224:C:H2'	1:XA:225:C:C6	2.44	0.53
8:XH:121:ASP:HB2	8:XH:125:ARG:NH2	2.24	0.53
11:XK:84:VAL:HG11	11:XK:95:ILE:HD11	1.91	0.53
25:YA:860:U:H5	25:YA:917:A:C2	2.25	0.53
29:YF:29:ASN:CB	29:YF:112:MET:HE1	2.38	0.53
30:YG:67:LYS:HE2	50:Y4:6:HIS:CE1	2.43	0.53
31:YH:40:GLU:O	31:YH:41:MET:HB2	2.09	0.53
36:YQ:81:VAL:C	36:YQ:82:ARG:CG	2.76	0.53
39:YT:16:ARG:HE	39:YT:19:LEU:HD21	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:YT:105:LEU:O	39:YT:107:ASP:N	2.42	0.53
53:Y7:9:ARG:NH1	53:Y7:47:ARG:HG3	2.24	0.53
4:QD:22:LYS:HG3	4:QD:26:CYS:SG	2.49	0.53
19:QS:44:MET:O	19:QS:46:GLY:N	2.40	0.53
25:RA:1681:G:HO2'	25:RA:1762:A:HO2'	1.49	0.53
25:RA:1803:A:O2'	27:RD:259:THR:HG21	2.09	0.53
28:RE:14:ILE:HG23	28:RE:15:PHE:N	2.22	0.53
30:RG:3:LEU:CD2	50:R4:25:TYR:CE1	2.91	0.53
45:RZ:163:LEU:H	45:RZ:163:LEU:HD12	1.72	0.53
47:R1:80:LEU:HD23	47:R1:80:LEU:H	1.74	0.53
48:R2:65:ASN:HB3	48:R2:69:ARG:NH2	2.24	0.53
50:R4:54:GLY:O	50:R4:71:ARG:HA	2.08	0.53
51:R5:55:ARG:HG3	51:R5:57:VAL:H	1.74	0.53
1:XA:769:G:H4'	1:XA:1513:A:H4'	1.90	0.53
18:XR:25:THR:HB	18:XR:26:LEU:HD23	1.91	0.53
25:YA:26:G:H1'	25:YA:515:A:N6	2.24	0.53
25:YA:1870:C:H2'	25:YA:1871:A:O4'	2.08	0.53
25:YA:1918:A:O2'	25:YA:1920:C:N4	2.42	0.53
26:YB:44:G:O2'	26:YB:47:C:N4	2.42	0.53
26:YB:52:A:N6	38:YS:33:LYS:HG3	2.24	0.53
27:YD:25:THR:HG21	27:YD:81:ALA:CB	2.38	0.53
27:YD:77:ALA:HB2	27:YD:97:TYR:CG	2.44	0.53
27:YD:233:HIS:CD2	27:YD:233:HIS:N	2.75	0.53
31:YH:12:PRO:HG3	31:YH:48:GLY:O	2.09	0.53
31:YH:139:GLN:O	31:YH:143:GLN:HB2	2.09	0.53
36:YQ:76:LYS:O	36:YQ:88:GLY:HA3	2.09	0.53
38:YS:10:ARG:O	38:YS:14:VAL:HG12	2.09	0.53
40:YU:102:GLU:OE1	41:YV:13:ARG:NH2	2.42	0.53
48:Y2:16:LEU:O	48:Y2:16:LEU:CG	2.49	0.53
1:QA:1371:G:C6	1:QA:1372:U:C4	2.97	0.53
3:QC:161:GLU:OE2	3:QC:161:GLU:HA	2.09	0.53
12:QL:27:LEU:C	12:QL:29:GLY:N	2.64	0.53
25:RA:1184:G:OP1	49:R3:29:ARG:NH1	2.41	0.53
25:RA:1614:A:N1	42:RW:91:GLY:HA2	2.24	0.53
25:RA:2250:G:C4	36:RQ:82:ARG:HG3	2.44	0.53
25:RA:2816:C:H2'	25:RA:2817:G:H8	1.74	0.53
28:RE:20:ALA:O	28:RE:21:VAL:CG2	2.48	0.53
30:RG:65:GLY:O	50:R4:7:PRO:HG2	2.09	0.53
31:RH:40:GLU:O	31:RH:41:MET:HB2	2.08	0.53
48:R2:4:SER:OG	48:R2:5:GLU:OE2	2.16	0.53
50:R4:15:ILE:HD13	50:R4:15:ILE:H	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:1004:A:H8	1:XA:1036:G:N2	2.05	0.53
1:XA:1128:C:H5'	9:XI:16:ARG:HH22	1.73	0.53
20:XT:95:ALA:O	20:XT:97:ALA:N	2.42	0.53
27:YD:25:THR:O	27:YD:26:LYS:C	2.52	0.53
27:YD:35:LYS:HD3	27:YD:63:ARG:CA	2.39	0.53
28:YE:119:ARG:HD3	28:YE:160:TYR:HB2	1.91	0.53
29:YF:9:ILE:HD11	29:YF:125:LEU:CG	2.36	0.53
29:YF:116:ASP:OD1	29:YF:119:ARG:NH2	2.41	0.53
32:YI:133:HIS:HB2	32:YI:134:PRO:HD2	1.91	0.53
33:YN:13:TRP:O	33:YN:135:PRO:HD2	2.08	0.53
33:YN:96:GLU:HG2	33:YN:97:ARG:N	2.24	0.53
48:Y2:50:ILE:CD1	48:Y2:51:ARG:N	2.61	0.53
52:Y6:26:ASN:ND2	52:Y6:35:GLU:OE2	2.42	0.53
1:QA:244:U:OP2	17:QQ:100:LYS:NZ	2.42	0.53
25:RA:224:G:O6	25:RA:419:C:O2'	2.27	0.53
25:RA:2512:C:H4'	28:RE:122:PHE:CE2	2.44	0.53
31:RH:10:PRO:C	31:RH:11:VAL:HG22	2.34	0.53
32:RI:133:HIS:HB2	32:RI:134:PRO:CD	2.39	0.53
36:RQ:76:LYS:O	36:RQ:88:GLY:HA3	2.09	0.53
42:RW:110:LYS:HG3	42:RW:111:HIS:ND1	2.23	0.53
54:R8:43:GLN:C	54:R8:44:LYS:HD2	2.34	0.53
25:YA:2784:C:H5"	28:YE:41:LYS:HZ3	1.74	0.53
31:YH:89:ILE:O	31:YH:89:ILE:CG1	2.57	0.53
32:YI:78:THR:HG22	32:YI:141:LYS:HG3	1.90	0.53
38:YS:56:LEU:C	38:YS:56:LEU:HD23	2.34	0.53
38:YS:56:LEU:O	38:YS:58:LEU:HD22	2.09	0.53
1:QA:1326:C:OP1	21:QU:17:THR:OG1	2.24	0.52
6:QF:10:LEU:HD13	6:QF:61:LEU:HD13	1.90	0.52
25:RA:1047:G:H2'	25:RA:1110:G:N1	2.24	0.52
25:RA:2115:G:N1	25:RA:2164:C:OP2	2.42	0.52
25:RA:2280:G:O2'	25:RA:2388:A:N1	2.38	0.52
25:RA:2327:A:H2'	25:RA:2328:A:C8	2.43	0.52
25:RA:2753:A:O2'	55:R9:15:LYS:NZ	2.39	0.52
28:RE:64:LYS:C	28:RE:66:HIS:N	2.68	0.52
29:RF:32:LEU:O	29:RF:36:VAL:HG23	2.09	0.52
30:RG:3:LEU:CD1	50:R4:25:TYR:CZ	2.90	0.52
31:RH:126:PRO:HD2	31:RH:127:GLU:H	1.72	0.52
32:RI:40:THR:HG22	32:RI:42:SER:H	1.73	0.52
36:RQ:7:MET:HE1	36:RQ:73:PRO:HG3	1.91	0.52
1:XA:89:U:O2'	1:XA:90:C:OP1	2.27	0.52
1:XA:1347:G:N2	1:XA:1374:A:OP2	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:XD:22:LYS:NZ	4:XD:22:LYS:HB3	2.23	0.52
19:XS:31:ILE:HG23	19:XS:49:ILE:HA	1.91	0.52
25:YA:1217:C:OP1	40:YU:15:LYS:HE3	2.09	0.52
28:YE:39:PRO:HG2	28:YE:40:GLU:OE1	2.09	0.52
28:YE:64:LYS:C	28:YE:66:HIS:N	2.68	0.52
29:YF:129:PHE:O	29:YF:142:TRP:CD1	2.62	0.52
29:YF:140:LEU:O	29:YF:143:ALA:HB3	2.09	0.52
30:YG:96:ARG:O	30:YG:98:ARG:N	2.42	0.52
31:YH:10:PRO:C	31:YH:11:VAL:HG22	2.34	0.52
31:YH:44:VAL:O	31:YH:44:VAL:CG2	2.57	0.52
44:YY:35:TYR:CE2	44:YY:69:ALA:HB3	2.44	0.52
50:Y4:18:CYS:SG	50:Y4:39:CYS:HB3	2.46	0.52
52:Y6:40:CYS:HB2	52:Y6:45:LYS:HD3	1.90	0.52
1:QA:658:G:OP1	15:QO:8:LYS:NZ	2.41	0.52
1:QA:1095:U:OP1	1:QA:1108:G:N2	2.42	0.52
13:QM:66:LEU:HA	13:QM:70:LEU:HB2	1.91	0.52
25:RA:1140:C:OP2	33:RN:66:LYS:NZ	2.42	0.52
28:RE:7:VAL:O	28:RE:196:VAL:HG13	2.09	0.52
28:RE:61:ARG:O	28:RE:63:LEU:N	2.42	0.52
31:RH:12:PRO:HG3	31:RH:48:GLY:O	2.09	0.52
31:RH:44:VAL:O	31:RH:44:VAL:CG2	2.57	0.52
50:R4:47:GLN:O	50:R4:48:ARG:HB2	2.07	0.52
2:XB:44:LEU:H	2:XB:44:LEU:HD12	1.74	0.52
7:XG:15:ASP:HB3	7:XG:19:GLY:H	1.74	0.52
12:XL:46:LYS:HG2	12:XL:47:LYS:H	1.73	0.52
25:YA:784:A:N7	27:YD:229:VAL:CG2	2.64	0.52
25:YA:1454:U:H5'	37:YR:63:ARG:HE	1.73	0.52
25:YA:2758:A:C4	31:YH:67:LEU:HD21	2.43	0.52
26:YB:66:A:O2'	26:YB:67:G:OP2	2.24	0.52
28:YE:7:VAL:O	28:YE:196:VAL:HG13	2.09	0.52
31:YH:121:ILE:HG12	31:YH:135:GLY:HA3	1.91	0.52
33:YN:110:GLY:O	33:YN:114:ARG:HG3	2.09	0.52
10:QJ:50:ILE:HD11	10:QJ:57:LYS:CG	2.35	0.52
25:RA:1655:A:O3'	28:RE:115:GLY:HA3	2.09	0.52
27:RD:35:LYS:NZ	27:RD:64:ILE:O	2.41	0.52
27:RD:206:LEU:O	27:RD:211:ARG:NH1	2.38	0.52
32:RI:86:THR:H	32:RI:123:LEU:HD12	1.74	0.52
35:RP:19:VAL:HG12	35:RP:27:HIS:HB2	1.91	0.52
41:RV:60:GLU:HB2	41:RV:97:LYS:HE3	1.92	0.52
50:R4:48:ARG:CZ	50:R4:51:ASP:HA	2.40	0.52
1:XA:191:G:O2'	20:XT:101:GLY:O	2.27	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:757:U:H2'	1:XA:758:G:O4'	2.10	0.52
25:YA:607:U:H3	25:YA:621:A:H2	1.52	0.52
25:YA:862:G:H2'	25:YA:863:A:O4'	2.09	0.52
25:YA:1311:G:N2	25:YA:1603:A:H62	2.07	0.52
25:YA:1438:U:H2'	25:YA:1439:A:C8	2.45	0.52
25:YA:2635:C:OP1	28:YE:78:LEU:HD12	2.09	0.52
25:YA:2666:C:N4	31:YH:109:PHE:HA	2.23	0.52
26:YB:48:A:H4'	38:YS:95:HIS:HD2	1.75	0.52
26:YB:48:A:H4'	38:YS:95:HIS:CD2	2.44	0.52
27:YD:155:LEU:CD1	27:YD:155:LEU:N	2.71	0.52
29:YF:196:LEU:C	29:YF:197:ASP:O	2.50	0.52
36:YQ:29:PHE:HB3	36:YQ:65:PHE:CZ	2.44	0.52
38:YS:25:ARG:CB	38:YS:25:ARG:HH11	2.22	0.52
54:Y8:61:LEU:O	54:Y8:62:LEU:CB	2.57	0.52
1:QA:544:G:OP2	4:QD:66:ARG:NH2	2.42	0.52
1:QA:920:U:H2'	1:QA:921:U:C6	2.44	0.52
1:QA:1292:U:H2'	1:QA:1293:G:C8	2.44	0.52
1:QA:1292:U:H2'	1:QA:1293:G:H8	1.73	0.52
1:QA:1492:A:OP1	12:QL:47:LYS:HB3	2.10	0.52
12:QL:47:LYS:HA	12:QL:48:PRO:C	2.35	0.52
25:RA:1854:A:H62	25:RA:1888:G:H8	1.57	0.52
25:RA:2557:G:H2'	25:RA:2558:C:H6	1.75	0.52
30:RG:82:LEU:HA	30:RG:86:MET:SD	2.48	0.52
35:RP:13:ASN:C	35:RP:15:ARG:H	2.17	0.52
36:RQ:29:PHE:HB3	36:RQ:65:PHE:CZ	2.44	0.52
50:R4:49:PHE:CD1	50:R4:49:PHE:N	2.77	0.52
1:XA:1336:C:O2'	1:XA:1337:G:O5'	2.25	0.52
28:YE:7:VAL:CG2	28:YE:8:LYS:H	2.10	0.52
28:YE:137:HIS:HB3	28:YE:138:PRO:CD	2.37	0.52
29:YF:132:VAL:O	29:YF:133:ASN:C	2.52	0.52
29:YF:162:LEU:HD23	29:YF:165:ARG:NH2	2.25	0.52
29:YF:192:LEU:HD21	29:YF:194:MET:CE	2.35	0.52
1:QA:1142:G:H3'	1:QA:1143:G:H8	1.75	0.52
1:QA:1224:G:C6	1:QA:1322:C:H1'	2.45	0.52
2:QB:134:GLU:HA	2:QB:137:ARG:HB3	1.92	0.52
13:QM:8:GLU:OE1	30:RG:115:ARG:CD	2.57	0.52
13:QM:121:LYS:HE2	13:QM:121:LYS:CA	2.36	0.52
25:RA:1028:A:N6	25:RA:1125:G:H2'	2.24	0.52
27:RD:133:LEU:HB3	27:RD:173:VAL:HG11	1.91	0.52
37:RR:59:ASP:OD2	37:RR:61:HIS:HB3	2.08	0.52
1:XA:56:U:H2'	1:XA:57:G:C8	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:1318:A:H5"	1:XA:1319:A:OP2	2.10	0.52
3:XC:40:ARG:O	3:XC:44:GLU:HB2	2.09	0.52
10:XJ:78:ASN:O	10:XJ:81:THR:OG1	2.25	0.52
25:YA:185:U:H4'	25:YA:218:A:H4'	1.92	0.52
25:YA:2089:U:H5"	25:YA:2089:U:C6	2.44	0.52
27:YD:66:ASP:OD2	27:YD:69:ARG:HG2	2.09	0.52
28:YE:93:VAL:C	28:YE:95:ILE:H	2.17	0.52
31:YH:24:VAL:HG23	31:YH:24:VAL:O	2.09	0.52
32:YI:115:ALA:HB3	32:YI:128:LEU:HD12	1.92	0.52
38:YS:62:LYS:HB3	38:YS:97:ARG:CD	2.39	0.52
48:Y2:7:ARG:HG3	48:Y2:7:ARG:NH1	2.25	0.52
3:QC:37:GLN:NE2	14:QN:52:GLN:OE1	2.32	0.52
25:RA:884:C:O2	25:RA:892:G:C2	2.62	0.52
25:RA:1043:C:N3	25:RA:1112:G:N2	2.46	0.52
25:RA:2776:A:OP1	25:RA:2776:A:H3'	2.10	0.52
28:RE:170:LEU:CD2	28:RE:185:LYS:HB2	2.40	0.52
36:RQ:81:VAL:HG23	46:R0:7:LEU:HD13	1.91	0.52
50:R4:55:ARG:C	50:R4:59:PHE:HB3	2.35	0.52
1:XA:593:G:H1	1:XA:646:U:H3	1.57	0.52
1:XA:964:A:O2'	10:XJ:55:LYS:NZ	2.40	0.52
3:XC:189:ALA:HB3	3:XC:196:LEU:HB2	1.91	0.52
25:YA:67:U:C4	25:YA:74:A:N1	2.78	0.52
25:YA:573:G:O2'	25:YA:574:C:H3'	2.09	0.52
25:YA:1790:C:H5"	25:YA:1791:A:OP1	2.09	0.52
25:YA:2233:U:H2'	25:YA:2234:G:C8	2.44	0.52
25:YA:2419:U:O4	54:Y8:30:ARG:NE	2.42	0.52
27:YD:133:LEU:HG	27:YD:189:CYS:O	2.10	0.52
28:YE:10:GLY:HA3	39:YT:8:LYS:HD2	1.91	0.52
29:YF:108:LYS:O	29:YF:112:MET:HG3	2.10	0.52
1:QA:64:G:H4'	1:QA:65:U:O5'	2.10	0.52
1:QA:115:G:H4'	1:QA:116:A:O5'	2.10	0.52
1:QA:1080:A:H5"	5:QE:16:THR:HG21	1.90	0.52
1:QA:1298:C:H4'	1:QA:1299:A:C4	2.45	0.52
12:QL:32:PHE:HE1	12:QL:86:ARG:HG3	1.73	0.52
25:RA:623:G:H2'	25:RA:624:C:C6	2.45	0.52
25:RA:2557:G:H2'	25:RA:2558:C:C6	2.44	0.52
27:RD:108:PRO:HB3	27:RD:143:HIS:HE1	1.73	0.52
27:RD:206:LEU:HD22	27:RD:211:ARG:HG2	1.92	0.52
28:RE:39:PRO:HG2	28:RE:40:GLU:OE1	2.09	0.52
31:RH:121:ILE:HG12	31:RH:135:GLY:HA3	1.91	0.52
44:RY:84:ARG:O	44:RY:95:LYS:HD3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:RY:87:LYS:HA	44:RY:92:ASN:HB3	1.91	0.52
50:R4:40:HIS:N	50:R4:41:PRO:CD	2.73	0.52
1:XA:64:G:H4'	1:XA:65:U:O5'	2.10	0.52
2:XB:35:GLU:O	2:XB:36:ARG:HD3	2.10	0.52
16:XP:8:ARG:O	16:XP:9:PHE:HD1	1.93	0.52
27:YD:35:LYS:HG2	27:YD:64:ILE:HG22	1.92	0.52
27:YD:36:PRO:HA	27:YD:62:TYR:O	2.09	0.52
28:YE:176:ILE:HG22	28:YE:176:ILE:O	2.10	0.52
29:YF:29:ASN:H	29:YF:112:MET:HE3	1.75	0.52
38:YS:56:LEU:O	38:YS:57:LYS:C	2.53	0.52
52:Y6:13:CYS:O	52:Y6:21:TYR:HA	2.09	0.52
1:QA:368:U:OP1	32:YI:91:SER:OG	2.27	0.52
1:QA:1372:U:H2'	1:QA:1373:G:O4'	2.10	0.52
1:QA:1426:C:H42	1:QA:1474:G:H1	1.58	0.52
25:RA:747:U:C5	51:R5:2:ALA:O	2.63	0.52
25:RA:1525:G:H2'	25:RA:1526:G:C8	2.45	0.52
25:RA:2335:A:O2'	25:RA:2336:A:O5'	2.25	0.52
25:RA:2867:G:O2'	25:RA:2868:A:H8	1.92	0.52
28:RE:116:VAL:HG22	28:RE:122:PHE:HB2	1.91	0.52
28:RE:179:GLU:OE1	28:RE:179:GLU:HA	2.10	0.52
31:RH:4:ILE:O	31:RH:6:ARG:N	2.43	0.52
31:RH:24:VAL:HG23	31:RH:24:VAL:O	2.09	0.52
41:RV:34:GLU:O	41:RV:36:PRO:HD3	2.10	0.52
50:R4:36:CYS:O	50:R4:37:SER:C	2.52	0.52
1:XA:1014:A:C2	1:XA:1219:U:H1'	2.45	0.52
7:XG:78:ARG:HG3	7:XG:79:ARG:N	2.25	0.52
19:XS:19:VAL:HG11	19:XS:44:MET:HG2	1.91	0.52
25:YA:2121:G:O6	25:YA:2176:A:N6	2.43	0.52
25:YA:2691:C:O3'	25:YA:2871:C:H4'	2.10	0.52
27:YD:174:ILE:N	27:YD:174:ILE:CD1	2.73	0.52
28:YE:15:PHE:CE1	39:YT:81:PRO:CD	2.93	0.52
38:YS:67:ARG:HB2	38:YS:67:ARG:HH11	1.65	0.52
38:YS:89:ARG:HG2	38:YS:89:ARG:HH11	1.74	0.52
38:YS:106:ARG:HA	38:YS:110:LEU:CG	2.39	0.52
51:Y5:56:LYS:HD3	51:Y5:58:LEU:HD23	1.92	0.52
52:Y6:7:ILE:HG13	52:Y6:8:LYS:H	1.75	0.52
54:Y8:43:GLN:C	54:Y8:44:LYS:HD2	2.35	0.52
1:QA:560:U:H5'	1:QA:566:G:N2	2.24	0.52
12:QL:127:GLU:O	12:QL:128:ALA:HB3	2.10	0.52
25:RA:395:U:H2'	25:RA:396:G:N7	2.24	0.52
25:RA:1178:C:H2'	25:RA:1179:C:C6	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1203:G:H3'	25:RA:1204:A:H5"	1.92	0.52
25:RA:1935:G:H1'	25:RA:1964:G:N2	2.24	0.52
26:RB:15:A:H5'	26:RB:16:G:C8	2.45	0.52
28:RE:129:HIS:O	28:RE:130:GLY:C	2.53	0.52
50:R4:14:ILE:HG23	50:R4:14:ILE:O	2.10	0.52
1:XA:964:A:N3	1:XA:969:A:O2'	2.37	0.52
2:XB:170:GLU:O	2:XB:174:VAL:HG23	2.10	0.52
25:YA:2470:G:H5'	36:YQ:56:ARG:HH22	1.75	0.52
25:YA:2593:U:H2'	25:YA:2594:C:C6	2.45	0.52
28:YE:61:ARG:O	28:YE:63:LEU:N	2.43	0.52
28:YE:170:LEU:CD2	28:YE:185:LYS:HB2	2.40	0.52
30:YG:88:ILE:O	30:YG:88:ILE:HD13	2.09	0.52
42:YW:40:ASN:O	42:YW:41:LYS:HG2	2.10	0.52
1:QA:300:A:O2'	1:QA:564:C:N3	2.40	0.52
1:QA:401:C:O2'	1:QA:621:A:N3	2.42	0.52
1:QA:1435:G:H2'	1:QA:1436:U:C6	2.44	0.52
11:QK:32:ILE:HG13	11:QK:72:ALA:HB2	1.92	0.52
28:RE:54:GLN:NE2	28:RE:54:GLN:H	2.08	0.52
30:RG:67:LYS:NZ	50:R4:6:HIS:CD2	2.78	0.52
31:RH:41:MET:HE1	31:RH:64:LEU:HB3	1.91	0.52
37:RR:33:ARG:HD3	37:RR:113:LEU:HG	1.92	0.52
42:RW:60:ASN:HD22	42:RW:60:ASN:H	1.56	0.52
43:RX:59:VAL:HG21	43:RX:78:LYS:HE3	1.91	0.52
50:R4:54:GLY:HA2	50:R4:57:GLU:HG2	1.92	0.52
1:XA:719:C:O2'	18:XR:50:ILE:O	2.24	0.52
1:XA:1007:C:H42	1:XA:1022:G:H1	1.56	0.52
1:XA:1491:G:C5	58:XA:1715:PAR:H21	2.45	0.52
2:XB:29:ALA:O	2:XB:32:ILE:HG22	2.10	0.52
12:XL:23:LYS:O	12:XL:24:VAL:HG23	2.10	0.52
14:XN:43:CYS:HA	14:XN:46:GLU:HG3	1.92	0.52
22:XV:16:C:N4	22:XV:17:U:O4	2.43	0.52
25:YA:1071:G:O6	25:YA:1091:G:O6	2.28	0.52
25:YA:1568:G:O4'	27:YD:58:HIS:HE1	1.91	0.52
25:YA:1932:A:H2'	25:YA:1933:G:O4'	2.09	0.52
25:YA:2760:C:H2'	25:YA:2761:G:H5"	1.91	0.52
25:YA:2795:G:H3'	25:YA:2797:U:C5'	2.40	0.52
27:YD:43:ARG:NH1	27:YD:44:ASN:OD1	2.42	0.52
27:YD:165:ILE:C	27:YD:166:GLN:HE21	2.18	0.52
27:YD:205:VAL:O	27:YD:206:LEU:C	2.52	0.52
31:YH:153:LYS:HG3	31:YH:161:GLY:HA2	1.91	0.52
36:YQ:119:ARG:HG2	36:YQ:119:ARG:NH1	2.20	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:YT:88:ILE:HD12	39:YT:90:GLN:N	2.25	0.52
45:YZ:89:PHE:HE2	45:YZ:96:VAL:HG21	1.75	0.52
48:Y2:15:LYS:H	48:Y2:67:LYS:HE2	1.73	0.52
49:Y3:40:THR:HB	49:Y3:43:ILE:HG12	1.92	0.52
51:Y5:49:CYS:SG	51:Y5:60:VAL:HG13	2.49	0.52
1:QA:501:C:O3'	12:QL:118:SER:HB2	2.10	0.51
1:QA:1064:G:HO2'	1:QA:1065:U:P	2.33	0.51
8:QH:77:GLU:HG2	8:QH:78:GLN:H	1.74	0.51
10:QJ:31:GLY:HA3	10:QJ:78:ASN:ND2	2.25	0.51
25:RA:270(F):U:H2'	25:RA:270(G):C:C6	2.45	0.51
25:RA:1190:G:H5'	35:RP:32:THR:HA	1.92	0.51
25:RA:1516:U:H2'	25:RA:1517:G:C8	2.44	0.51
25:RA:2277:G:C5'	36:RQ:85:LYS:HG3	2.39	0.51
28:RE:54:GLN:O	28:RE:55:ASN:HB2	2.09	0.51
28:RE:93:VAL:C	28:RE:95:ILE:H	2.17	0.51
31:RH:55:PRO:HG2	31:RH:61:HIS:ND1	2.25	0.51
35:RP:111:ARG:HG2	35:RP:128:HIS:CD2	2.45	0.51
47:R1:2:SER:HB2	47:R1:4:VAL:HG12	1.92	0.51
1:XA:1432:G:OP1	39:YT:107:ASP:HB2	2.10	0.51
1:XA:1525:G:P	11:XK:120:ARG:HH22	2.32	0.51
4:XD:13:ARG:HD2	4:XD:38:TYR:O	2.10	0.51
4:XD:108:LEU:HB3	4:XD:110:PHE:CE1	2.45	0.51
12:XL:127:GLU:O	12:XL:128:ALA:HB3	2.10	0.51
25:YA:2056:G:H2'	25:YA:2056:G:N3	2.25	0.51
25:YA:2246:G:H2'	25:YA:2247:A:C8	2.45	0.51
25:YA:2287:A:N6	25:YA:2344:U:C2	2.77	0.51
28:YE:54:GLN:O	28:YE:55:ASN:HB2	2.09	0.51
28:YE:95:ILE:H	28:YE:95:ILE:CD1	2.19	0.51
29:YF:125:LEU:HA	29:YF:194:MET:O	2.10	0.51
30:YG:77:ILE:HD13	30:YG:82:LEU:HD12	1.92	0.51
47:Y1:83:GLU:HG2	47:Y1:84:GLY:N	2.24	0.51
25:RA:1055:G:H1	25:RA:1104:C:H42	1.57	0.51
27:RD:12:SER:HB2	27:RD:208:LYS:HB3	1.92	0.51
50:R4:23:GLU:C	50:R4:24:THR:HG1	2.18	0.51
51:R5:40:LYS:HD3	51:R5:46:CYS:SG	2.50	0.51
1:XA:632:A:C8	1:XA:633:G:C8	2.97	0.51
25:YA:190:A:OP2	47:Y1:39:LYS:HE3	2.10	0.51
25:YA:620:G:H4'	25:YA:621:A:C5'	2.40	0.51
25:YA:712:G:H1	25:YA:719:C:N4	2.08	0.51
25:YA:1190:G:H5'	35:YP:32:THR:HA	1.92	0.51
26:YB:12:C:O2'	46:Y0:74:ARG:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:67:PHE:CE2	27:YD:157:ARG:NH2	2.79	0.51
41:YV:25:LEU:H	41:YV:92:THR:HG21	1.74	0.51
18:QR:32:ARG:HA	18:QR:69:THR:HG21	1.91	0.51
25:RA:49:A:C8	25:RA:120:U:C5	2.98	0.51
25:RA:483:A:H5'	44:RY:49:VAL:HG22	1.93	0.51
25:RA:674:G:H1'	29:RF:74:ARG:HD3	1.92	0.51
25:RA:1204:A:H1'	25:RA:1206:G:C4	2.46	0.51
25:RA:1728:G:H3'	25:RA:1729:A:H5"	1.91	0.51
25:RA:2712:U:O2'	25:RA:2712(A):A:C8	2.62	0.51
25:RA:2818:G:OP2	37:RR:42:LYS:NZ	2.43	0.51
28:RE:14:ILE:CD1	39:RT:14:TYR:CZ	2.93	0.51
28:RE:105:THR:HB	28:RE:197:ILE:HG12	1.92	0.51
28:RE:105:THR:HG23	28:RE:166:THR:OG1	2.10	0.51
30:RG:5:VAL:CG2	50:R4:25:TYR:CE2	2.94	0.51
31:RH:89:ILE:O	31:RH:89:ILE:CG1	2.57	0.51
31:RH:126:PRO:HD2	31:RH:127:GLU:N	2.26	0.51
36:RQ:64:ILE:HA	36:RQ:106:VAL:CG1	2.33	0.51
51:R5:20:ARG:C	51:R5:22:HIS:N	2.68	0.51
3:XC:47:LEU:HD11	3:XC:76:VAL:HB	1.91	0.51
5:XE:33:VAL:HG11	5:XE:109:ILE:HA	1.91	0.51
9:XI:11:LYS:H	9:XI:104:ARG:HH21	1.57	0.51
22:XV:19:G:H5'	22:XV:20:U:H5	1.76	0.51
25:YA:307:G:H21	25:YA:330:A:H62	1.57	0.51
25:YA:754:C:H2'	25:YA:755:C:C6	2.46	0.51
25:YA:2298:A:H62	25:YA:2318:G:H8	1.58	0.51
27:YD:30:GLU:HG3	27:YD:63:ARG:NH2	2.26	0.51
28:YE:203:LYS:HE3	28:YE:204:ALA:CB	2.40	0.51
30:YG:105:LYS:HD3	50:Y4:26:SER:HB3	1.93	0.51
37:YR:83:ILE:HG22	37:YR:87:TYR:HE2	1.76	0.51
40:YU:107:ALA:O	40:YU:110:VAL:HB	2.10	0.51
1:QA:1269:A:N1	1:QA:1312:G:O2'	2.34	0.51
5:QE:71:LEU:HD22	5:QE:115:VAL:HG22	1.93	0.51
25:RA:1101:U:H2'	25:RA:1102:C:H6	1.76	0.51
30:RG:6:ALA:HB2	50:R4:23:GLU:CG	2.40	0.51
35:RP:59:LEU:HD12	35:RP:61:ARG:NH1	2.25	0.51
42:RW:67:ASP:OD1	42:RW:67:ASP:N	2.33	0.51
43:RX:35:THR:HG23	43:RX:38:GLU:HG2	1.93	0.51
44:RY:74:PRO:O	44:RY:80:GLY:HA2	2.11	0.51
1:XA:302:G:O3'	12:XL:17:LYS:HE2	2.11	0.51
1:XA:401:C:H2'	1:XA:402:G:C8	2.45	0.51
1:XA:598:U:H4'	8:XH:94:TYR:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:991:U:O4	1:XA:1212:U:O2'	2.18	0.51
5:XE:13:ILE:HD11	5:XE:55:VAL:HG22	1.91	0.51
9:XI:70:LYS:O	9:XI:74:ILE:HG13	2.10	0.51
12:XL:24:VAL:O	12:XL:24:VAL:CG1	2.58	0.51
25:YA:43:G:H2'	25:YA:44:A:O4'	2.11	0.51
25:YA:451:C:H4'	29:YF:52:LYS:HZ1	1.76	0.51
25:YA:751:A:H5'	42:YW:90:ARG:HA	1.92	0.51
25:YA:957:A:N1	25:YA:2458:G:H4'	2.26	0.51
25:YA:1291:C:H5'	25:YA:1536:A:H5'	1.92	0.51
27:YD:35:LYS:HD2	27:YD:104:TYR:CD1	2.45	0.51
33:YN:38:HIS:O	40:YU:67:ALA:HB1	2.10	0.51
51:Y5:42:PRO:HB2	51:Y5:43:HIS:ND1	2.24	0.51
1:QA:503:C:OP2	12:QL:116:SER:HB3	2.11	0.51
1:QA:1220:G:O3'	19:QS:36:ARG:HD3	2.11	0.51
21:QU:6:ARG:HE	21:QU:15:ARG:HH21	1.58	0.51
25:RA:1266:G:O5'	42:RW:15:ARG:NH2	2.42	0.51
25:RA:1354:A:OP1	27:RD:38:LYS:HE2	2.10	0.51
25:RA:1636:C:H2'	25:RA:1637:A:C8	2.46	0.51
25:RA:2663:G:H3'	25:RA:2664:G:H8	1.75	0.51
36:RQ:25:ASP:HA	36:RQ:100:GLY:O	2.11	0.51
20:XT:10:LEU:O	20:XT:13:LEU:HG	2.11	0.51
25:YA:296:C:N3	25:YA:342:G:O6	2.44	0.51
25:YA:443:A:H1'	25:YA:1201:C:O4'	2.09	0.51
27:YD:221:VAL:HG22	27:YD:226:MET:HE2	1.91	0.51
28:YE:20:ALA:C	28:YE:21:VAL:HG13	2.35	0.51
28:YE:51:PHE:CD2	28:YE:52:LEU:N	2.76	0.51
29:YF:65:TRP:HZ2	29:YF:72:ARG:NH2	2.09	0.51
32:YI:124:GLY:H	32:YI:142:VAL:HG23	1.75	0.51
39:YT:109:GLU:O	39:YT:113:LYS:HB2	2.11	0.51
1:QA:564:C:C5	17:QQ:31:LEU:HD11	2.46	0.51
1:QA:625:G:H2'	1:QA:626:U:C6	2.46	0.51
3:QC:79:ARG:CZ	11:XK:99:GLN:HB2	2.41	0.51
4:QD:31:CYS:N	4:QD:32:ALA:HA	2.25	0.51
20:QT:14:LYS:HA	20:QT:17:ARG:HG3	1.91	0.51
25:RA:49:A:H61	25:RA:177:G:H2'	1.76	0.51
25:RA:2646:C:H2'	25:RA:2647:U:O4'	2.11	0.51
27:RD:43:ARG:HH11	27:RD:44:ASN:CG	2.16	0.51
28:RE:20:ALA:C	28:RE:21:VAL:HG13	2.35	0.51
28:RE:119:ARG:HD3	28:RE:160:TYR:HB2	1.91	0.51
39:RT:28:VAL:HG23	39:RT:88:ILE:HA	1.92	0.51
45:RZ:182:LYS:H	45:RZ:182:LYS:HD3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:R4:12:ALA:HB1	50:R4:30:GLU:H	1.76	0.51
53:R7:22:MET:HE2	53:R7:31:LEU:CD1	2.40	0.51
1:XA:266:G:H5'	1:XA:266:G:N9	2.25	0.51
1:XA:836:G:C6	1:XA:851:G:C6	2.98	0.51
4:XD:78:LEU:HD22	4:XD:96:LEU:HB3	1.92	0.51
6:XF:4:TYR:HD1	6:XF:92:LYS:HA	1.76	0.51
7:XG:155:ARG:NH2	7:XG:155:ARG:O	2.43	0.51
13:XM:122:LYS:O	13:XM:122:LYS:HD3	2.11	0.51
17:XQ:100:LYS:O	17:XQ:101:ARG:NE	2.43	0.51
19:XS:15:LEU:O	19:XS:19:VAL:N	2.36	0.51
25:YA:952:G:P	36:YQ:16:ARG:HH12	2.34	0.51
25:YA:2283:C:H2'	25:YA:2284:C:O4'	2.10	0.51
30:YG:16:ARG:O	30:YG:20:ILE:HG12	2.10	0.51
31:YH:41:MET:HE1	31:YH:64:LEU:HB3	1.91	0.51
41:YV:65:GLY:HA3	41:YV:91:TYR:CZ	2.46	0.51
45:YZ:10:ARG:NH2	45:YZ:26:GLY:O	2.44	0.51
47:Y1:70:VAL:O	47:Y1:74:VAL:HG23	2.10	0.51
51:Y5:38:ALA:HB3	51:Y5:40:LYS:HE3	1.92	0.51
54:Y8:29:LYS:HB2	54:Y8:44:LYS:HG2	1.90	0.51
1:QA:719:C:O2'	18:QR:49:LYS:HB3	2.10	0.51
1:QA:1252:A:H61	1:QA:1285:A:H61	1.58	0.51
1:QA:1318:A:H4'	19:QS:11:VAL:CG1	2.40	0.51
22:QV:21:A:H61	22:QV:46:G:H2'	1.73	0.51
25:RA:747:U:C5	51:R5:2:ALA:C	2.89	0.51
36:RQ:83:MET:HB2	46:R0:7:LEU:HD12	1.91	0.51
38:RS:56:LEU:HD23	38:RS:58:LEU:HD22	1.92	0.51
54:R8:10:ALA:O	54:R8:14:VAL:HG12	2.11	0.51
54:R8:61:LEU:O	54:R8:62:LEU:CB	2.57	0.51
1:XA:779:C:H2'	1:XA:780:A:O4'	2.11	0.51
1:XA:1002:G:H1	1:XA:1038:C:N4	2.06	0.51
6:XF:97:PHE:CD1	18:XR:31:LEU:HD21	2.46	0.51
13:XM:14:ARG:N	13:XM:44:ARG:HD3	2.21	0.51
19:XS:26:GLY:O	19:XS:28:LYS:N	2.43	0.51
25:YA:1914:C:H2'	25:YA:1915:U:O4'	2.11	0.51
25:YA:2747:G:O6	25:YA:2755:C:H5"	2.10	0.51
27:YD:210:GLY:O	27:YD:213:ARG:N	2.43	0.51
38:YS:95:HIS:CG	38:YS:96:GLY:N	2.77	0.51
42:YW:106:ILE:O	42:YW:106:ILE:HG12	2.07	0.51
44:YY:81:LYS:HG2	44:YY:97:ARG:HD3	1.93	0.51
49:Y3:43:ILE:O	49:Y3:47:VAL:HG23	2.10	0.51
54:Y8:52:LYS:N	54:Y8:53:PRO:HD2	2.22	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:660:G:N2	1:QA:745:C:N3	2.53	0.51
1:QA:1004:A:H1'	1:QA:1036:G:N2	2.26	0.51
1:QA:1356:G:H2'	1:QA:1357:A:C8	2.46	0.51
19:QS:69:HIS:CE1	50:R4:69:LYS:HE2	2.44	0.51
25:RA:507:A:H5"	25:RA:508:G:H5'	1.91	0.51
25:RA:1930:G:O2'	25:RA:1931:U:P	2.69	0.51
25:RA:2037:G:H2'	25:RA:2038:G:C8	2.45	0.51
28:RE:47:VAL:O	28:RE:48:GLN:C	2.52	0.51
31:RH:131:VAL:CG1	31:RH:132:ARG:N	2.74	0.51
40:RU:90:VAL:HG22	41:RV:39:LEU:HB3	1.93	0.51
1:XA:67:C:H2'	1:XA:68:G:H8	1.76	0.51
1:XA:1002:G:H2'	1:XA:1003:G:C8	2.45	0.51
1:XA:1075:C:OP1	2:XB:179:LYS:HE2	2.11	0.51
1:XA:1232:U:OP1	9:XI:124:GLN:NE2	2.43	0.51
2:XB:21:ARG:HB2	2:XB:39:ILE:HA	1.91	0.51
4:XD:30:LYS:C	4:XD:32:ALA:H	2.18	0.51
20:XT:89:ARG:HH21	20:XT:104:LEU:HD11	1.76	0.51
25:YA:321:G:C5	29:YF:165:ARG:NH1	2.79	0.51
25:YA:656:G:H2'	25:YA:657:U:O4'	2.11	0.51
25:YA:1021:A:H61	25:YA:1142(A):A:N6	2.09	0.51
27:YD:117:VAL:CG2	27:YD:128:GLY:C	2.84	0.51
27:YD:134:ARG:HB2	27:YD:135:PHE:HD1	1.75	0.51
31:YH:55:PRO:HG2	31:YH:61:HIS:ND1	2.26	0.51
31:YH:72:ILE:O	31:YH:75:ALA:HB3	2.11	0.51
36:YQ:25:ASP:HA	36:YQ:100:GLY:O	2.11	0.51
37:YR:104:ARG:HD3	37:YR:111:LEU:HD21	1.92	0.51
51:Y5:45:VAL:HG11	51:Y5:57:VAL:HG12	1.93	0.51
53:Y7:9:ARG:NH2	53:Y7:48:LYS:CB	2.62	0.51
1:QA:49:U:H5	1:QA:365:U:O4	1.91	0.51
1:QA:836:G:C6	1:QA:851:G:C6	2.99	0.51
25:RA:628:G:HO2'	25:RA:651:G:HO2'	1.55	0.51
25:RA:2725:A:O2'	25:RA:2726:U:H5"	2.10	0.51
28:RE:37:ARG:HE	28:RE:37:ARG:N	2.09	0.51
29:RF:135:LYS:HB3	29:RF:138:GLU:HG3	1.93	0.51
36:RQ:36:ALA:HB1	36:RQ:127:ILE:HD12	1.93	0.51
43:RX:40:LYS:HG3	43:RX:51:VAL:HB	1.92	0.51
50:R4:42:PHE:O	50:R4:44:THR:N	2.44	0.51
1:XA:1016:A:H2'	1:XA:1017:G:O4'	2.11	0.51
12:XL:126:LYS:C	12:XL:128:ALA:N	2.69	0.51
13:XM:36:LYS:C	13:XM:36:LYS:HD3	2.36	0.51
15:XO:33:THR:HG21	15:XO:85:LEU:HD22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:740:U:H2'	25:YA:741:G:C8	2.45	0.51
25:YA:2208:U:H1'	27:YD:151:LYS:HE2	1.92	0.51
27:YD:28:GLU:OE1	27:YD:29:PRO:HD2	2.11	0.51
28:YE:37:ARG:HE	28:YE:37:ARG:N	2.09	0.51
28:YE:54:GLN:NE2	28:YE:54:GLN:H	2.08	0.51
31:YH:19:VAL:HG13	31:YH:43:VAL:CG2	2.41	0.51
32:YI:62:LYS:HE3	32:YI:134:PRO:HG2	1.93	0.51
35:YP:61:ARG:CD	54:Y8:13:ARG:HD2	2.40	0.51
36:YQ:58:PHE:O	36:YQ:58:PHE:HD1	1.94	0.51
36:YQ:133:ARG:HG2	36:YQ:134:ARG:N	2.26	0.51
38:YS:83:LYS:HG2	38:YS:109:GLY:HA2	1.90	0.51
38:YS:86:ALA:O	38:YS:87:PHE:HB3	2.09	0.51
48:Y2:36:ARG:O	48:Y2:40:SER:HB2	2.10	0.51
1:QA:1321:C:H4'	13:QM:87:TYR:CZ	2.46	0.51
9:QI:8:GLY:HA2	9:QI:79:LEU:HD12	1.91	0.51
10:QJ:55:LYS:CG	10:QJ:56:HIS:N	2.70	0.51
12:QL:11:VAL:HG13	17:QQ:29:HIS:CD2	2.45	0.51
25:RA:1678:G:H22	25:RA:1989:G:H22	1.57	0.51
28:RE:137:HIS:HB3	28:RE:138:PRO:CD	2.37	0.51
30:RG:88:ILE:O	30:RG:88:ILE:HD13	2.10	0.51
31:RH:72:ILE:O	31:RH:75:ALA:HB3	2.11	0.51
36:RQ:23:GLY:O	36:RQ:24:GLY:C	2.54	0.51
39:RT:118:ARG:HH21	39:RT:121:ILE:HG21	1.76	0.51
50:R4:39:CYS:HB3	50:R4:41:PRO:CD	2.40	0.51
51:R5:36:CYS:C	51:R5:38:ALA:H	2.19	0.51
1:XA:390:C:O3'	16:XP:28:ARG:NH2	2.44	0.51
2:XB:73:THR:OG1	2:XB:170:GLU:OE2	2.23	0.51
11:XK:41:THR:HG21	11:XK:71:LYS:HB3	1.93	0.51
25:YA:2629:A:O2'	25:YA:2630:G:H5"	2.11	0.51
31:YH:103:LEU:CD1	31:YH:131:VAL:HG21	2.41	0.51
33:YN:42:TRP:O	40:YU:64:ARG:NH2	2.37	0.51
38:YS:74:ALA:O	38:YS:75:GLU:C	2.54	0.51
1:QA:359:U:H2'	1:QA:360:A:H8	1.74	0.50
1:QA:1065:U:O5'	1:QA:1190:G:N2	2.44	0.50
1:QA:1152:A:H5"	10:QJ:13:HIS:CD2	2.46	0.50
1:QA:1237:C:H5"	1:QA:1238:A:O4'	2.11	0.50
8:QH:20:TYR:HE2	8:QH:75:ARG:HD2	1.76	0.50
12:QL:23:LYS:O	12:QL:24:VAL:HG23	2.10	0.50
16:QP:3:LYS:HG3	16:QP:24:ALA:HB2	1.92	0.50
19:QS:5:LEU:HD22	50:R4:67:TYR:CE2	2.45	0.50
25:RA:1268:A:H2'	25:RA:1269:A:O4'	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1833:U:O2'	25:RA:1969:A:N1	2.33	0.50
25:RA:2511:U:O4	25:RA:2575:C:N3	2.44	0.50
26:RB:17:C:H2'	26:RB:18:G:O4'	2.10	0.50
31:RH:153:LYS:HG3	31:RH:161:GLY:HA2	1.91	0.50
36:RQ:133:ARG:HG2	36:RQ:134:ARG:N	2.26	0.50
1:XA:1118:C:OP1	9:XI:9:ARG:HD3	2.11	0.50
1:XA:1358:U:OP1	14:XN:35:ARG:HG3	2.11	0.50
3:XC:138:VAL:HG22	3:XC:151:VAL:HG23	1.93	0.50
5:XE:94:ALA:HB2	5:XE:119:LEU:HG	1.92	0.50
25:YA:84:A:O5'	44:YY:8:LYS:HD3	2.11	0.50
25:YA:1068:G:O2'	25:YA:1096:A:H1'	2.11	0.50
25:YA:1550:C:H2'	25:YA:1551:C:C6	2.46	0.50
25:YA:2619:C:O2'	28:YE:156:MET:HE3	2.11	0.50
26:YB:37:C:O2	38:YS:95:HIS:NE2	2.44	0.50
31:YH:126:PRO:HD2	31:YH:127:GLU:N	2.25	0.50
36:YQ:7:MET:HE1	36:YQ:73:PRO:HG3	1.91	0.50
38:YS:83:LYS:HG2	38:YS:109:GLY:H	1.76	0.50
41:YV:52:VAL:HG23	41:YV:55:ALA:H	1.75	0.50
42:YW:57:ASN:O	42:YW:61:ASN:HB2	2.10	0.50
48:Y2:17:SER:CB	48:Y2:18:PRO:CA	2.89	0.50
1:QA:126:G:OP1	1:QA:605:U:O2'	2.23	0.50
12:QL:126:LYS:C	12:QL:128:ALA:N	2.69	0.50
25:RA:195:A:H61	25:RA:198:C:H3'	1.76	0.50
25:RA:844:C:H2'	25:RA:845:G:O4'	2.11	0.50
25:RA:1416:G:H2'	25:RA:1417:C:H6	1.76	0.50
28:RE:37:ARG:HE	28:RE:37:ARG:H	1.59	0.50
32:RI:1:MET:N	32:RI:21:VAL:O	2.42	0.50
33:RN:46:VAL:HG13	33:RN:48:MET:HG3	1.93	0.50
39:RT:19:LEU:HD22	39:RT:86:ILE:HG22	1.94	0.50
1:XA:947:G:H2'	1:XA:948:C:C6	2.46	0.50
1:XA:1120:G:H2'	1:XA:1121:U:C6	2.46	0.50
1:XA:1315:U:H2'	1:XA:1316:G:O4'	2.11	0.50
1:XA:1318:A:O2'	19:XS:37:ARG:HD3	2.11	0.50
2:XB:9:GLU:HB3	2:XB:48:MET:SD	2.50	0.50
2:XB:80:ILE:HD11	2:XB:208:ILE:HG23	1.93	0.50
5:XE:67:VAL:HG21	5:XE:140:ARG:HA	1.92	0.50
6:XF:86:ARG:O	6:XF:87:ARG:HG2	2.11	0.50
10:XJ:50:ILE:HD11	10:XJ:57:LYS:HD3	1.93	0.50
10:XJ:57:LYS:CD	10:XJ:60:ARG:CZ	2.89	0.50
19:XS:4:SER:O	19:XS:5:LEU:HD13	2.11	0.50
19:XS:42:PRO:HB3	50:Y4:60:GLN:OE1	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:XT:26:ASN:HB3	20:XT:71:THR:OG1	2.10	0.50
25:YA:27:G:N2	25:YA:512:G:H2'	2.25	0.50
25:YA:507:A:H5"	25:YA:508:G:H5'	1.93	0.50
25:YA:587:C:OP2	35:YP:21:ARG:NH2	2.44	0.50
28:YE:116:VAL:HG22	28:YE:122:PHE:HB2	1.91	0.50
28:YE:129:HIS:O	28:YE:130:GLY:C	2.53	0.50
31:YH:4:ILE:O	31:YH:6:ARG:N	2.43	0.50
38:YS:26:LEU:CD2	38:YS:87:PHE:CD1	2.94	0.50
38:YS:87:PHE:O	38:YS:88:ASP:O	2.29	0.50
38:YS:89:ARG:HD2	38:YS:89:ARG:O	2.11	0.50
1:QA:1152:A:H2'	1:QA:1153:C:C6	2.46	0.50
7:QG:116:ALA:O	7:QG:120:ILE:HG12	2.11	0.50
9:QI:53:VAL:HB	9:QI:95:LYS:HE3	1.92	0.50
25:RA:2844:G:H3'	25:RA:2845:G:H8	1.75	0.50
26:RB:15:A:H1'	26:RB:109:G:N9	2.26	0.50
51:R5:37:LYS:HD2	51:R5:37:LYS:O	2.12	0.50
1:XA:64:G:H4'	1:XA:65:U:C5'	2.42	0.50
1:XA:1032(A):G:H2'	1:XA:1032(B):G:C8	2.47	0.50
3:XC:54:ARG:HD3	3:XC:56:ASP:OD1	2.10	0.50
20:XT:83:ARG:O	20:XT:86:ARG:HB3	2.11	0.50
25:YA:336:C:HO2'	44:YY:35:TYR:HH	1.55	0.50
25:YA:1844:C:H2'	25:YA:1845:G:H8	1.76	0.50
27:YD:10:THR:HG23	27:YD:13:ARG:CB	2.34	0.50
27:YD:233:HIS:CD2	27:YD:233:HIS:H	2.29	0.50
28:YE:105:THR:HG23	28:YE:166:THR:OG1	2.10	0.50
29:YF:32:LEU:O	29:YF:36:VAL:HG23	2.11	0.50
29:YF:108:LYS:HA	29:YF:108:LYS:NZ	2.27	0.50
30:YG:5:VAL:HG11	30:YG:100:TRP:HB3	1.93	0.50
1:QA:19:C:P	5:QE:127:ASN:HD22	2.35	0.50
1:QA:790:A:OP1	22:QV:38:A:O2'	2.25	0.50
11:QK:17:GLY:N	11:QK:79:SER:O	2.44	0.50
25:RA:705:A:H1'	27:RD:9:TYR:CE2	2.46	0.50
25:RA:784:A:O2'	25:RA:785:G:H5"	2.12	0.50
25:RA:998:C:H2'	25:RA:999:U:O4'	2.12	0.50
31:RH:16:SER:O	31:RH:17:VAL:HG23	2.12	0.50
31:RH:169:VAL:HG13	31:RH:170:ARG:N	2.26	0.50
32:RI:120:ILE:HD11	32:RI:126:TYR:CZ	2.46	0.50
42:RW:60:ASN:HD22	42:RW:60:ASN:N	2.09	0.50
1:XA:309:G:H2'	1:XA:310:G:H8	1.77	0.50
2:XB:21:ARG:O	2:XB:23:ARG:N	2.43	0.50
2:XB:114:ARG:O	2:XB:117:GLU:HB2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:XO:6:GLU:H	15:XO:6:GLU:CD	2.15	0.50
19:XS:40:ILE:HG23	19:XS:67:VAL:O	2.12	0.50
25:YA:1268:A:H2'	25:YA:1269:A:O4'	2.12	0.50
30:YG:101:ILE:CD1	50:Y4:9:LEU:HD11	2.41	0.50
38:YS:26:LEU:HD22	38:YS:87:PHE:CD1	2.46	0.50
50:Y4:10:VAL:HG22	50:Y4:11:PRO:HD2	1.94	0.50
10:QJ:47:PHE:CE1	10:QJ:63:PHE:HB2	2.47	0.50
10:QJ:51:ARG:NH2	14:QN:58:LYS:HZ1	2.10	0.50
11:QK:33:THR:HG22	11:QK:39:PRO:HA	1.92	0.50
13:QM:33:ALA:HA	13:QM:59:TYR:HE2	1.76	0.50
13:QM:40:ASN:ND2	13:QM:43:THR:HG23	2.27	0.50
25:RA:2378:A:O5'	25:RA:2378:A:H8	1.93	0.50
27:RD:62:TYR:CE2	27:RD:64:ILE:HA	2.46	0.50
28:RE:61:ARG:CB	28:RE:62:PRO:HD3	2.41	0.50
28:RE:176:ILE:HG22	28:RE:176:ILE:O	2.10	0.50
28:RE:203:LYS:HE3	28:RE:204:ALA:CB	2.40	0.50
29:RF:157:VAL:HG21	29:RF:181:LEU:HD21	1.93	0.50
35:RP:36:LYS:HD2	35:RP:37:GLY:H	1.76	0.50
44:RY:97:ARG:HH21	44:RY:98:VAL:HB	1.77	0.50
47:R1:62:VAL:HG23	47:R1:63:ALA:O	2.11	0.50
51:R5:54:GLY:O	51:R5:55:ARG:C	2.54	0.50
1:XA:1392:G:N2	1:XA:1502:A:H8	2.09	0.50
2:XB:111:ARG:HH21	2:XB:114:ARG:HG2	1.76	0.50
25:YA:796:C:H2'	25:YA:797:C:C6	2.46	0.50
25:YA:2133:G:H1'	25:YA:2158:A:N6	2.26	0.50
25:YA:2469:A:H5'	25:YA:2470:G:OP2	2.11	0.50
27:YD:182:LEU:H	27:YD:272:ALA:CB	2.25	0.50
31:YH:131:VAL:CG1	31:YH:132:ARG:N	2.74	0.50
31:YH:169:VAL:HG13	31:YH:170:ARG:N	2.26	0.50
32:YI:33:ARG:HB3	32:YI:35:LEU:HG	1.92	0.50
34:YO:120:GLU:HG2	34:YO:122:LEU:HG	1.94	0.50
36:YQ:108:GLY:O	36:YQ:109:VAL:HG23	2.12	0.50
36:YQ:132:VAL:HG12	36:YQ:133:ARG:N	2.27	0.50
48:Y2:41:ILE:HD11	48:Y2:44:LEU:CB	2.42	0.50
56:Z6:101:PPU:HN2	56:Z6:101:PPU:HD2	1.76	0.50
1:QA:826:C:H2'	1:QA:827:U:O2	2.11	0.50
2:QB:231:GLU:HG3	2:QB:233:SER:H	1.77	0.50
9:QI:95:LYS:NZ	9:QI:96:LEU:HD13	2.26	0.50
13:QM:40:ASN:HD22	13:QM:43:THR:HG23	1.77	0.50
13:QM:121:LYS:HZ3	24:QY:40:G:H4'	1.77	0.50
25:RA:650:C:OP1	54:R8:48:PHE:CZ	2.64	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1007:C:H5”	33:RN:35:ARG:NH1	2.27	0.50
28:RE:14:ILE:CG1	39:RT:14:TYR:CZ	2.95	0.50
31:RH:133:VAL:HG12	31:RH:141:VAL:HG13	1.93	0.50
32:RI:94:ALA:H	32:RI:116:LEU:HD13	1.77	0.50
36:RQ:108:GLY:O	36:RQ:109:VAL:HG23	2.12	0.50
39:RT:39:ARG:HG2	39:RT:40:THR:H	1.76	0.50
50:R4:1:MET:HE3	50:R4:6:HIS:CD2	2.46	0.50
51:R5:48:GLU:HA	51:R5:59:GLU:HG2	1.94	0.50
56:Z5:101:PPU:HN2	56:Z5:101:PPU:HD2	1.76	0.50
1:XA:407:G:H2’	1:XA:408:A:C8	2.47	0.50
1:XA:1145:C:H4’	1:XA:1146:A:H5’	1.92	0.50
1:XA:1414:U:H2’	1:XA:1415:G:C8	2.46	0.50
2:XB:189:ASP:HB3	2:XB:203:GLY:O	2.12	0.50
4:XD:108:LEU:HD21	4:XD:183:GLY:HA3	1.93	0.50
11:XK:86:GLY:O	11:XK:91:ARG:HD3	2.11	0.50
27:YD:241:PRO:O	27:YD:242:ARG:C	2.53	0.50
31:YH:133:VAL:HG12	31:YH:141:VAL:HG13	1.93	0.50
41:YV:61:VAL:HA	41:YV:94:LEU:HD23	1.93	0.50
43:YX:6:ASP:OD2	48:Y2:29:LYS:NZ	2.45	0.50
48:Y2:9:GLN:O	48:Y2:12:GLU:HB3	2.10	0.50
1:QA:701:C:H1’	1:QA:703:G:C5	2.46	0.50
2:QB:162:ILE:HD11	2:QB:184:VAL:HG22	1.93	0.50
4:QD:22:LYS:HE2	4:QD:31:CYS:O	2.11	0.50
9:QI:118:LYS:O	9:QI:120:ARG:N	2.40	0.50
25:RA:270(L):U:H2’	32:RI:50:ARG:HD2	1.93	0.50
25:RA:516:C:OP1	51:R5:13:LYS:NZ	2.45	0.50
25:RA:1300:U:H4’	25:RA:1301:A:H5”	1.93	0.50
25:RA:2543:G:H2’	25:RA:2544:G:C8	2.46	0.50
31:RH:19:VAL:HG13	31:RH:43:VAL:HG23	1.93	0.50
35:RP:88:LEU:HD12	35:RP:95:VAL:HG11	1.93	0.50
38:RS:67:ARG:O	38:RS:71:ARG:HG3	2.12	0.50
39:RT:111:ARG:O	39:RT:113:LYS:N	2.42	0.50
51:R5:50:GLY:O	51:R5:51:TYR:CB	2.59	0.50
53:R7:5:TRP:NE1	53:R7:7:PRO:HG3	2.26	0.50
1:XA:245:C:C2	1:XA:284:G:C2	3.00	0.50
4:XD:112:VAL:HG12	4:XD:116:GLN:OE1	2.12	0.50
25:YA:2291:U:H2’	25:YA:2292:C:C6	2.46	0.50
28:YE:2:LYS:HG2	28:YE:95:ILE:CG2	2.42	0.50
28:YE:105:THR:HB	28:YE:197:ILE:HG12	1.92	0.50
31:YH:24:VAL:HG21	31:YH:72:ILE:HG12	1.94	0.50
36:YQ:36:ALA:HB1	36:YQ:127:ILE:HD12	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:Y7:5:TRP:NE1	53:Y7:7:PRO:HG3	2.26	0.50
1:QA:1172:C:H2'	1:QA:1173:G:H8	1.75	0.50
25:RA:1285:G:N2	25:RA:1328:G:H5"	2.27	0.50
25:RA:1496:A:H8	25:RA:1577:C:O2'	1.94	0.50
25:RA:1588:C:H2'	25:RA:1589:C:C6	2.47	0.50
25:RA:2469:A:H5'	25:RA:2470:G:OP2	2.12	0.50
28:RE:2:LYS:HG2	28:RE:95:ILE:CG2	2.42	0.50
36:RQ:2:LEU:HD23	36:RQ:2:LEU:N	2.27	0.50
43:RX:26:TYR:HB3	43:RX:92:LEU:HD12	1.93	0.50
50:R4:1:MET:HG3	50:R4:1:MET:O	2.12	0.50
1:XA:532:A:O2'	1:XA:533:A:OP1	2.27	0.50
1:XA:677:U:H3	1:XA:713:G:H22	1.60	0.50
1:XA:767:A:H2'	1:XA:768:A:O4'	2.11	0.50
1:XA:826:C:H2'	1:XA:827:U:O2	2.11	0.50
1:XA:1399:C:C2	1:XA:1401:G:C5	3.00	0.50
2:XB:32:ILE:HD11	2:XB:40:HIS:HB3	1.94	0.50
15:XO:67:LEU:HB3	15:XO:78:TYR:HE1	1.77	0.50
25:YA:528:A:N1	25:YA:2042:A:H2'	2.26	0.50
25:YA:1847:A:H8	25:YA:1847:A:OP1	1.95	0.50
27:YD:35:LYS:HE2	27:YD:104:TYR:HB2	1.94	0.50
28:YE:9:VAL:HG11	39:YT:7:ILE:CG2	2.42	0.50
31:YH:16:SER:O	31:YH:17:VAL:HG23	2.12	0.50
33:YN:58:ASP:OD1	33:YN:58:ASP:N	2.45	0.50
38:YS:35:ILE:CD1	38:YS:101:LEU:HD23	2.41	0.50
44:YY:97:ARG:HH21	44:YY:98:VAL:HB	1.76	0.50
52:Y6:47:THR:HG22	52:Y6:48:VAL:HG12	1.94	0.50
1:QA:1073:U:OP2	5:QE:57:LYS:NZ	2.35	0.50
4:QD:28:SER:HB3	4:QD:29:PRO:HD2	1.94	0.50
13:QM:89:GLY:O	13:QM:92:HIS:HB2	2.11	0.50
19:QS:42:PRO:CG	50:R4:63:TYR:HE2	2.24	0.50
20:QT:64:ASP:CG	20:QT:81:LYS:HZ2	2.19	0.50
25:RA:669:G:H2'	25:RA:669:G:N3	2.27	0.50
25:RA:848:G:H2'	25:RA:849:A:C8	2.47	0.50
25:RA:1525:G:H2'	25:RA:1526:G:H8	1.76	0.50
25:RA:1639:U:H2'	25:RA:1640:C:H5"	1.93	0.50
26:RB:38:C:N4	26:RB:44:G:H1	2.06	0.50
31:RH:19:VAL:HG13	31:RH:43:VAL:CG2	2.41	0.50
31:RH:103:LEU:CD1	31:RH:131:VAL:HG21	2.41	0.50
31:RH:153:LYS:O	31:RH:154:PRO:O	2.30	0.50
33:RN:34:LEU:O	33:RN:49:GLY:HA3	2.12	0.50
39:RT:34:VAL:HG12	39:RT:36:GLU:HG2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:RT:42:ILE:HD12	39:RT:42:ILE:H	1.77	0.50
50:R4:10:VAL:HG23	50:R4:11:PRO:HD2	1.93	0.50
50:R4:57:GLU:O	50:R4:61:ARG:O	2.30	0.50
50:R4:68:ARG:HD3	50:R4:69:LYS:HG2	1.92	0.50
1:XA:280:C:N3	17:XQ:39:SER:N	2.47	0.50
1:XA:606:G:H1	1:XA:631:G:H5"	1.77	0.50
1:XA:1126:U:H1'	1:XA:1280:A:C5	2.47	0.50
2:XB:162:ILE:HD11	2:XB:184:VAL:HG22	1.94	0.50
5:XE:76:ILE:HG13	5:XE:93:PRO:HB3	1.94	0.50
13:XM:49:THR:HB	13:XM:52:GLU:H	1.77	0.50
13:XM:121:LYS:NZ	24:XY:40:G:H4'	2.27	0.50
25:YA:363:G:H2'	25:YA:363(A):A:H8	1.77	0.50
25:YA:577:G:O2'	25:YA:1254:A:OP1	2.28	0.50
25:YA:1045:A:O2'	25:YA:1046:A:OP2	2.29	0.50
25:YA:1608:A:H1'	25:YA:1610:A:OP2	2.12	0.50
25:YA:2585:U:H5	56:Z6:101:PPU:HO2'	1.57	0.50
28:YE:14:ILE:HD11	39:YT:14:TYR:HH	1.72	0.50
31:YH:7:LEU:C	31:YH:7:LEU:HD12	2.37	0.50
36:YQ:23:GLY:O	36:YQ:24:GLY:C	2.54	0.50
54:Y8:40:GLU:C	54:Y8:42:ARG:N	2.68	0.50
1:QA:1278:U:O4	10:QJ:99:LYS:NZ	2.34	0.49
3:QC:14:ILE:HG12	3:QC:15:THR:N	2.27	0.49
8:QH:9:MET:HG3	8:QH:26:VAL:HG21	1.94	0.49
12:QL:24:VAL:O	12:QL:24:VAL:CG1	2.58	0.49
25:RA:1105:U:H2'	25:RA:1106:G:H8	1.77	0.49
25:RA:1637:A:H4'	25:RA:2711:A:O2'	2.11	0.49
25:RA:2745:C:H1'	31:RH:143:GLN:HG2	1.94	0.49
25:RA:2758:A:C2	25:RA:2759:G:H1'	2.47	0.49
28:RE:119:ARG:HD3	28:RE:160:TYR:CD2	2.47	0.49
32:RI:8:PRO:HD3	32:RI:15:VAL:HG13	1.94	0.49
32:RI:57:ARG:O	32:RI:61:ARG:HG2	2.12	0.49
36:RQ:80:GLU:HG3	36:RQ:81:VAL:N	2.27	0.49
39:RT:36:GLU:HG3	39:RT:41:ARG:CD	2.39	0.49
43:RX:40:LYS:O	43:RX:42:ALA:N	2.45	0.49
1:XA:299:G:H2'	1:XA:300:A:C8	2.47	0.49
1:XA:397:A:H5'	1:XA:398:C:OP1	2.12	0.49
1:XA:1356:G:H2'	1:XA:1357:A:C8	2.47	0.49
11:XK:82:VAL:HB	11:XK:108:ILE:HG12	1.94	0.49
19:XS:41:VAL:HB	19:XS:42:PRO:CA	2.42	0.49
20:XT:64:ASP:HA	20:XT:67:ALA:HB3	1.93	0.49
25:YA:273(C):C:N4	25:YA:363(C):G:H1	2.09	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1113:U:H2'	25:YA:1114:G:H8	1.77	0.49
25:YA:1210:A:H5'	25:YA:1210:A:C8	2.46	0.49
27:YD:2:ALA:CB	27:YD:20:ASP:CB	2.90	0.49
27:YD:2:ALA:CB	27:YD:20:ASP:HB3	2.42	0.49
27:YD:218:ARG:HB3	27:YD:219:PRO:HD2	1.94	0.49
28:YE:46:ALA:HB1	28:YE:80:GLU:HB2	1.94	0.49
29:YF:11:VAL:CG1	29:YF:12:LEU:N	2.75	0.49
29:YF:53:THR:C	29:YF:55:GLY:N	2.68	0.49
32:YI:120:ILE:HD11	32:YI:126:TYR:CZ	2.47	0.49
45:YZ:152:ALA:O	45:YZ:154:ASP:N	2.41	0.49
1:QA:921:U:O2	5:QE:19:MET:HB2	2.12	0.49
1:QA:1213:A:N6	1:QA:1215:G:N3	2.60	0.49
2:QB:235:SER:OG	2:QB:236:TYR:N	2.46	0.49
6:QF:97:PHE:O	18:QR:31:LEU:HD23	2.12	0.49
19:QS:41:VAL:HA	19:QS:44:MET:HG3	1.93	0.49
25:RA:140:A:H8	25:RA:1408:C:HO2'	1.60	0.49
25:RA:242:G:O5'	54:R8:3:LYS:HE3	2.11	0.49
25:RA:911:A:C5	36:RQ:9:TYR:CD2	3.00	0.49
25:RA:954:G:O2'	25:RA:2274:A:N1	2.38	0.49
25:RA:1209:G:O2'	25:RA:1237:A:N1	2.46	0.49
25:RA:1694:C:H4'	25:RA:1695:G:O5'	2.11	0.49
25:RA:1799:G:H4'	25:RA:1800:C:O5'	2.12	0.49
29:RF:178:PRO:HB2	29:RF:201:VAL:HG11	1.93	0.49
37:RR:44:LEU:HD22	37:RR:48:VAL:HG23	1.94	0.49
1:XA:337:C:H2'	1:XA:338:A:C8	2.48	0.49
1:XA:719:C:H1'	18:XR:49:LYS:HB3	1.93	0.49
1:XA:735:C:H5'	18:XR:71:LYS:HD3	1.95	0.49
1:XA:1281:U:H5'	1:XA:1282:C:OP2	2.11	0.49
16:XP:28:ARG:NH1	16:XP:29:ASP:OD2	2.45	0.49
20:XT:29:LYS:HD2	20:XT:66:ALA:CB	2.42	0.49
25:YA:1693:U:H1'	27:YD:14:ARG:NH2	2.27	0.49
25:YA:2235:G:H2'	25:YA:2236:C:C6	2.47	0.49
27:YD:76:PRO:HA	27:YD:118:VAL:HG23	1.93	0.49
38:YS:5:THR:HG1	38:YS:7:TYR:HB3	1.75	0.49
46:Y0:70:GLN:OE1	46:Y0:72:ARG:HD3	2.12	0.49
48:Y2:69:ARG:CB	48:Y2:69:ARG:HH11	2.25	0.49
1:QA:414:A:OP2	1:QA:428:G:N2	2.28	0.49
1:QA:501:C:OP1	12:QL:117:ARG:NH2	2.37	0.49
1:QA:1142:G:H2'	1:QA:1143:G:O4'	2.12	0.49
1:QA:1298:C:O2'	1:QA:1299:A:OP2	2.23	0.49
9:QI:46:ALA:HB2	9:QI:74:ILE:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:QV:75:C:H5”	22:QV:76:A:OP2	2.11	0.49
25:RA:1991:U:H2’	25:RA:1992:G:H5”	1.93	0.49
25:RA:2563:U:H1’	25:RA:2566:A:N6	2.27	0.49
25:RA:2689:U:H4’	25:RA:2690:C:O5’	2.12	0.49
25:RA:2734:A:H5’	25:RA:2735:G:OP2	2.12	0.49
25:RA:2811:G:OP1	28:RE:61:ARG:HG3	2.12	0.49
25:RA:2867:G:O2’	25:RA:2868:A:P	2.70	0.49
26:RB:24:G:H1’	26:RB:27:C:N4	2.27	0.49
30:RG:3:LEU:CD2	50:R4:25:TYR:CD1	2.87	0.49
30:RG:115:ARG:NH2	30:RG:137:GLU:OE1	2.45	0.49
31:RH:120:GLY:HA3	31:RH:140:LYS:NZ	2.28	0.49
45:RZ:166:SER:HB2	45:RZ:168:GLU:N	2.27	0.49
1:XA:255:G:O6	1:XA:266:G:O6	2.30	0.49
1:XA:464:G:C6	1:XA:466:C:H5’	2.48	0.49
1:XA:838:G:N1	1:XA:848:C:N4	2.32	0.49
13:XM:121:LYS:CA	13:XM:121:LYS:HE2	2.42	0.49
14:XN:23:ARG:NH1	14:XN:30:ALA:HB2	2.27	0.49
25:YA:918:A:N3	26:YB:80:U:O2’	2.43	0.49
25:YA:2832:U:H4’	25:YA:2833:G:H5”	1.93	0.49
28:YE:37:ARG:HE	28:YE:37:ARG:H	1.59	0.49
28:YE:61:ARG:CB	28:YE:62:PRO:CD	2.90	0.49
38:YS:52:SER:O	38:YS:56:LEU:CD2	2.60	0.49
1:QA:1327:C:H2’	1:QA:1328:C:H6	1.76	0.49
3:QC:14:ILE:HG12	3:QC:15:THR:H	1.76	0.49
3:QC:157:ILE:HD11	3:QC:166:GLU:HB2	1.95	0.49
8:QH:95:VAL:HB	8:QH:99:GLU:O	2.13	0.49
25:RA:594:U:H5’	54:R8:61:LEU:CD2	2.42	0.49
25:RA:724:U:H2’	25:RA:725:G:O4’	2.12	0.49
25:RA:1272:A:OP2	25:RA:1647:G:OP1	2.30	0.49
25:RA:1532:C:H42	25:RA:1539:G:H1	1.59	0.49
25:RA:2805:G:H2’	25:RA:2807:G:C8	2.48	0.49
27:RD:175:LEU:HD12	27:RD:185:VAL:HG21	1.93	0.49
28:RE:23:VAL:HG12	28:RE:173:VAL:HG21	1.94	0.49
28:RE:55:ASN:O	28:RE:57:LYS:N	2.44	0.49
31:RH:23:ARG:HD2	31:RH:34:GLU:OE2	2.12	0.49
36:RQ:90:VAL:C	36:RQ:92:GLY:N	2.70	0.49
38:RS:15:ARG:NH1	38:RS:25:ARG:HH21	2.11	0.49
50:R4:9:LEU:H	50:R4:27:THR:CG2	2.25	0.49
1:XA:724:G:C2	1:XA:725:G:C8	3.00	0.49
1:XA:974:A:OP2	14:XN:29:ARG:NH2	2.42	0.49
5:XE:11:ILE:CG2	5:XE:108:ALA:CB	2.90	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:XF:97:PHE:HB2	18:XR:32:ARG:CZ	2.41	0.49
25:YA:28:A:N6	25:YA:512:G:H1'	2.27	0.49
25:YA:242:G:H3'	54:Y8:6:THR:HG23	1.95	0.49
25:YA:613:U:H5'	25:YA:616:A:N6	2.28	0.49
25:YA:1586:A:H3'	25:YA:1587:A:H8	1.76	0.49
25:YA:2243:U:H2'	25:YA:2244:U:C6	2.48	0.49
25:YA:2319:G:N7	38:YS:3:ARG:HB3	2.27	0.49
25:YA:2635:C:H5'	28:YE:77:ILE:HD13	1.93	0.49
25:YA:2821:A:OP2	28:YE:110:GLY:HA3	2.12	0.49
27:YD:35:LYS:CG	27:YD:64:ILE:HG22	2.42	0.49
28:YE:17:ASP:N	28:YE:17:ASP:OD1	2.46	0.49
28:YE:18:ASP:O	28:YE:19:ARG:C	2.56	0.49
31:YH:153:LYS:O	31:YH:154:PRO:O	2.29	0.49
36:YQ:2:LEU:HD23	36:YQ:2:LEU:N	2.27	0.49
36:YQ:29:PHE:N	36:YQ:105:GLU:OE2	2.41	0.49
43:YX:53:LYS:HB3	43:YX:82:GLN:HB3	1.93	0.49
45:YZ:5:LEU:HB3	45:YZ:59:LEU:HA	1.94	0.49
50:Y4:15:ILE:H	50:Y4:15:ILE:HD13	1.76	0.49
54:Y8:10:ALA:O	54:Y8:14:VAL:HG12	2.11	0.49
1:QA:573:A:N3	1:QA:883:C:O2'	2.44	0.49
1:QA:1142:G:H3'	1:QA:1143:G:C8	2.48	0.49
14:QN:15:LYS:HD2	14:QN:16:PHE:CE2	2.47	0.49
18:QR:26:LEU:HD22	18:QR:42:ARG:HD2	1.94	0.49
25:RA:684:G:C2	25:RA:794:G:C2	3.00	0.49
25:RA:971:C:H2'	25:RA:972:G:O4'	2.12	0.49
25:RA:1257:C:H4'	29:RF:83:PHE:CD1	2.47	0.49
25:RA:2619:C:O2'	28:RE:156:MET:HE3	2.12	0.49
25:RA:2665:A:H2'	25:RA:2666:C:O4'	2.12	0.49
26:RB:78:A:H2'	26:RB:79:C:O4'	2.12	0.49
30:RG:110:ALA:HB1	30:RG:140:ILE:HD12	1.94	0.49
31:RH:103:LEU:H	31:RH:103:LEU:HD23	1.77	0.49
33:RN:7:LYS:H	33:RN:7:LYS:HZ3	1.59	0.49
36:RQ:29:PHE:N	36:RQ:105:GLU:OE2	2.41	0.49
39:RT:41:ARG:NH1	39:RT:41:ARG:HB2	2.26	0.49
40:RU:92:ARG:O	40:RU:92:ARG:HG2	2.12	0.49
43:RX:60:ARG:HH12	53:R7:47:ARG:HH22	1.59	0.49
54:R8:15:LYS:C	54:R8:15:LYS:HD3	2.37	0.49
1:XA:347:G:O2'	1:XA:348:G:H5"	2.12	0.49
1:XA:1129:C:O2'	1:XA:1131:G:N7	2.45	0.49
10:XJ:49:VAL:HG22	14:XN:41:ARG:HB2	1.94	0.49
13:XM:121:LYS:HE2	13:XM:121:LYS:N	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:994:C:OP1	40:YU:53:ARG:NH2	2.45	0.49
28:YE:111:ARG:HA	37:YR:1:MET:CG	2.43	0.49
28:YE:179:GLU:O	28:YE:180:ASN:HB2	2.12	0.49
29:YF:144:LYS:C	29:YF:146:ALA:H	2.21	0.49
30:YG:94:LEU:HD12	30:YG:99:MET:HA	1.95	0.49
31:YH:103:LEU:HD23	31:YH:103:LEU:H	1.77	0.49
33:YN:17:ASP:O	33:YN:56:ASN:HB2	2.12	0.49
33:YN:34:LEU:HD21	33:YN:120:LEU:HB2	1.94	0.49
38:YS:60:GLY:O	38:YS:61:ASN:CB	2.55	0.49
38:YS:89:ARG:O	38:YS:90:GLY:C	2.55	0.49
40:YU:95:LEU:HD22	41:YV:4:ILE:HD12	1.93	0.49
48:Y2:18:PRO:C	48:Y2:20:GLU:H	2.20	0.49
52:Y6:41:PRO:HD2	52:Y6:46:HIS:N	2.28	0.49
54:Y8:16:ILE:CD1	54:Y8:57:ARG:HG2	2.42	0.49
1:QA:109:A:C6	1:QA:326:G:C6	3.01	0.49
13:QM:7:VAL:HG21	30:RG:113:ARG:C	2.38	0.49
25:RA:242:G:C8	54:R8:5:LYS:HG2	2.46	0.49
25:RA:806:C:P	35:RP:41:ARG:HH11	2.35	0.49
25:RA:883:G:H22	25:RA:892:G:N2	2.10	0.49
25:RA:1022:G:O2'	25:RA:1023:U:OP2	2.20	0.49
25:RA:1278:A:H4'	37:RR:34:ILE:HD12	1.93	0.49
25:RA:2023:G:H5'	25:RA:2617:C:H4'	1.94	0.49
25:RA:2747:G:OP1	31:RH:138:LYS:NZ	2.45	0.49
30:RG:54:GLU:HA	30:RG:57:ALA:HB3	1.94	0.49
32:RI:144:VAL:HG22	32:RI:145:VAL:H	1.77	0.49
33:RN:134:ARG:N	33:RN:135:PRO:HD3	2.28	0.49
36:RQ:132:VAL:HG12	36:RQ:133:ARG:N	2.27	0.49
45:RZ:70:LEU:HB2	45:RZ:91:LEU:HD21	1.95	0.49
49:R3:6:VAL:HG13	49:R3:56:VAL:HG13	1.93	0.49
50:R4:42:PHE:O	50:R4:44:THR:O	2.31	0.49
53:R7:31:LEU:HD22	53:R7:42:LEU:HD13	1.95	0.49
1:XA:103:C:O2'	1:XA:172:A:N1	2.37	0.49
1:XA:108:G:H5"	1:XA:109:A:H5"	1.94	0.49
1:XA:729:A:H2'	1:XA:730:G:H8	1.76	0.49
1:XA:1225:A:H5"	1:XA:1226:C:OP2	2.13	0.49
2:XB:12:GLU:C	2:XB:14:GLY:H	2.21	0.49
13:XM:81:LEU:HD13	13:XM:88:ARG:HD2	1.94	0.49
25:YA:574:C:N3	28:YE:145:LYS:CE	2.76	0.49
25:YA:587:C:H4'	25:YA:588:U:O5'	2.12	0.49
25:YA:1478:G:H2'	25:YA:1479:G:H8	1.76	0.49
25:YA:2182:G:H2'	25:YA:2183:C:C6	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2820:A:O5'	37:YR:4:LEU:HD23	2.13	0.49
27:YD:30:GLU:CD	27:YD:63:ARG:HE	2.21	0.49
28:YE:179:GLU:OE1	28:YE:179:GLU:HA	2.10	0.49
38:YS:99:LYS:O	38:YS:101:LEU:N	2.45	0.49
40:YU:61:TRP:CD2	40:YU:94:ASN:HA	2.47	0.49
43:YX:57:LEU:HD11	43:YX:78:LYS:HD2	1.94	0.49
50:Y4:35:VAL:C	50:Y4:37:SER:H	2.20	0.49
52:Y6:15:GLU:CD	52:Y6:41:PRO:HB3	2.37	0.49
1:QA:579:G:H5'	1:QA:728:A:H1'	1.95	0.49
1:QA:584:G:H2'	1:QA:585:G:C8	2.48	0.49
3:QC:73:PRO:O	3:QC:76:VAL:HG22	2.13	0.49
17:QQ:18:THR:HG23	17:QQ:69:LYS:HE3	1.94	0.49
25:RA:26:G:H1'	25:RA:515:A:H61	1.77	0.49
25:RA:1085:A:O2'	25:RA:1086:A:OP1	2.25	0.49
25:RA:1667:G:H8	25:RA:1667:G:OP2	1.95	0.49
25:RA:1899:G:H21	25:RA:1902:C:H42	1.56	0.49
25:RA:1937:A:O2'	25:RA:1939:U:OP2	2.29	0.49
25:RA:2406:U:C2	35:RP:72:PRO:HB2	2.48	0.49
28:RE:95:ILE:H	28:RE:95:ILE:CD1	2.19	0.49
31:RH:54:ARG:HD3	31:RH:65:HIS:ND1	2.27	0.49
31:RH:98:LEU:HD12	31:RH:102:ALA:O	2.13	0.49
31:RH:124:GLU:HB3	31:RH:132:ARG:CG	2.42	0.49
31:RH:128:PRO:HD2	31:RH:129:THR:N	2.25	0.49
43:RX:27:THR:HB	43:RX:80:ILE:HB	1.94	0.49
45:RZ:152:ALA:O	45:RZ:154:ASP:N	2.41	0.49
50:R4:47:GLN:O	50:R4:48:ARG:CB	2.61	0.49
54:R8:33:ASN:O	54:R8:35:GLN:N	2.46	0.49
1:XA:401:C:H2'	1:XA:402:G:H8	1.78	0.49
1:XA:407:G:O4'	4:XD:119:GLN:NE2	2.46	0.49
1:XA:618:C:H5'	1:XA:619:U:H5"	1.95	0.49
1:XA:908:A:H2'	1:XA:909:A:C8	2.48	0.49
2:XB:24:TRP:CZ3	2:XB:26:PRO:HA	2.48	0.49
5:XE:10:MET:SD	5:XE:13:ILE:HD13	2.53	0.49
25:YA:729:G:O6	27:YD:209:ALA:N	2.44	0.49
25:YA:1055:G:O2'	25:YA:1085:A:N1	2.44	0.49
25:YA:1154:G:OP2	40:YU:58:ARG:NH1	2.46	0.49
29:YF:51:THR:O	29:YF:93:LYS:NZ	2.38	0.49
31:YH:19:VAL:HG13	31:YH:43:VAL:HG23	1.93	0.49
31:YH:54:ARG:HD3	31:YH:65:HIS:ND1	2.27	0.49
31:YH:98:LEU:HD12	31:YH:102:ALA:O	2.13	0.49
31:YH:124:GLU:HB3	31:YH:132:ARG:CG	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:YT:107:ASP:H	39:YT:110:ILE:HG22	1.78	0.49
44:YY:86:ARG:HB2	44:YY:95:LYS:HD2	1.94	0.49
46:Y0:53:MET:HA	46:Y0:58:THR:O	2.13	0.49
52:Y6:41:PRO:O	52:Y6:45:LYS:HE3	2.12	0.49
1:QA:31:G:O2'	1:QA:48:C:N4	2.45	0.49
1:QA:1015:A:H2'	1:QA:1016:A:C8	2.47	0.49
3:QC:47:LEU:HD23	3:QC:68:VAL:HG11	1.94	0.49
7:QG:20:ASP:HB3	7:QG:23:VAL:HG23	1.94	0.49
25:RA:607:U:OP1	29:RF:102:PRO:HA	2.13	0.49
25:RA:2212:A:H1'	25:RA:2215:G:C5	2.47	0.49
25:RA:2336:A:H61	46:R0:43:THR:CG2	2.25	0.49
25:RA:2393:A:H2'	25:RA:2394:C:O4'	2.13	0.49
29:RF:150:GLY:HA2	29:RF:172:TRP:CE3	2.48	0.49
29:RF:160:ASN:HB3	29:RF:163:VAL:HB	1.95	0.49
30:RG:60:LEU:O	30:RG:64:THR:HG22	2.11	0.49
31:RH:7:LEU:C	31:RH:7:LEU:HD12	2.37	0.49
36:RQ:23:GLY:O	36:RQ:24:GLY:O	2.30	0.49
44:RY:47:LYS:HG2	44:RY:60:PHE:HD1	1.77	0.49
50:R4:22:ILE:H	50:R4:22:ILE:HD12	1.77	0.49
52:R6:14:THR:O	52:R6:49:HIS:HA	2.12	0.49
1:XA:163:C:H2'	1:XA:164:U:C6	2.47	0.49
1:XA:474:G:H2'	1:XA:475:G:H8	1.77	0.49
1:XA:1079:G:C6	1:XA:1080:A:N6	2.81	0.49
3:XC:148:GLY:HA3	3:XC:172:ARG:O	2.12	0.49
6:XF:19:LEU:HD21	6:XF:59:TYR:CE1	2.47	0.49
13:XM:14:ARG:HG2	13:XM:17:VAL:HG23	1.95	0.49
25:YA:264:C:C2'	25:YA:265:A:H5"	2.43	0.49
25:YA:527:C:OP2	25:YA:2779:U:H5	1.96	0.49
25:YA:846:C:C2	25:YA:847:U:C5	3.00	0.49
25:YA:2277:G:H5"	36:YQ:85:LYS:CB	2.42	0.49
26:YB:12:C:C2'	46:Y0:73:GLY:HA3	2.43	0.49
34:YO:4:PRO:O	34:YO:5:GLN:HB2	2.11	0.49
38:YS:11:LYS:HG2	38:YS:11:LYS:O	2.12	0.49
52:Y6:21:TYR:HE2	52:Y6:53:LYS:HE3	1.77	0.49
54:Y8:58:ILE:O	54:Y8:61:LEU:HD12	2.13	0.49
1:QA:523:A:H61	12:QL:92:ASP:HB2	1.76	0.49
1:QA:564:C:OP1	12:QL:15:ARG:NE	2.45	0.49
1:QA:625:G:H2'	1:QA:626:U:H6	1.78	0.49
1:QA:976:G:N2	1:QA:1362(A):C:OP2	2.29	0.49
2:QB:21:ARG:O	2:QB:23:ARG:N	2.46	0.49
2:QB:96:ARG:H	2:QB:96:ARG:HD2	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:QS:68:GLY:C	50:R4:68:ARG:HG2	2.38	0.49
28:RE:46:ALA:HB1	28:RE:80:GLU:HB2	1.94	0.49
28:RE:119:ARG:HD3	28:RE:160:TYR:HD2	1.78	0.49
31:RH:12:PRO:HD3	31:RH:48:GLY:O	2.13	0.49
31:RH:24:VAL:HG21	31:RH:72:ILE:HG12	1.94	0.49
50:R4:10:VAL:CG2	50:R4:11:PRO:HD2	2.43	0.49
51:R5:2:ALA:O	51:R5:3:LYS:CB	2.61	0.49
51:R5:48:GLU:HA	51:R5:59:GLU:CG	2.43	0.49
54:R8:58:ILE:O	54:R8:61:LEU:HD12	2.13	0.49
1:XA:109:A:C6	1:XA:326:G:C6	3.01	0.49
1:XA:901:A:C5	1:XA:902:G:H1'	2.48	0.49
1:XA:1238:A:H62	1:XA:1301:U:H3	1.59	0.49
1:XA:1297:C:H1'	1:XA:1298:C:H5	1.78	0.49
2:XB:47:THR:HA	2:XB:202:PRO:HG2	1.95	0.49
10:XJ:35:SER:OG	10:XJ:73:ASP:HB2	2.13	0.49
12:XL:11:VAL:HG11	17:XQ:36:ILE:HG21	1.93	0.49
20:XT:89:ARG:NH2	20:XT:104:LEU:HD11	2.27	0.49
25:YA:639:U:H2'	25:YA:640:C:C6	2.48	0.49
25:YA:1266:G:OP2	51:Y5:20:ARG:NE	2.41	0.49
25:YA:2729:G:O2'	28:YE:187:ALA:HB2	2.12	0.49
27:YD:69:ARG:C	27:YD:71:ASP:N	2.69	0.49
27:YD:123:ALA:HB3	27:YD:131:LEU:HG	1.94	0.49
27:YD:227:ASN:HB3	27:YD:228:PRO:CD	2.30	0.49
28:YE:119:ARG:HD3	28:YE:160:TYR:CD2	2.47	0.49
35:YP:147:LEU:HB3	35:YP:148:LEU:H	1.41	0.49
36:YQ:34:LEU:HD23	36:YQ:104:PHE:HD2	1.77	0.49
54:Y8:15:LYS:HD3	54:Y8:15:LYS:C	2.37	0.49
1:QA:110:C:H2'	1:QA:111:G:O4'	2.13	0.49
1:QA:1510:U:H2'	1:QA:1511:G:C8	2.47	0.49
4:QD:9:CYS:SG	4:QD:22:LYS:CE	2.90	0.49
4:QD:20:TYR:HA	4:QD:26:CYS:SG	2.53	0.49
5:QE:76:ILE:CG2	5:QE:118:ILE:HD13	2.43	0.49
7:QG:155:ARG:O	7:QG:155:ARG:NH2	2.46	0.49
10:QJ:24:VAL:HG21	10:QJ:37:PRO:HD3	1.95	0.49
14:QN:27:CYS:SG	14:QN:29:ARG:HB2	2.53	0.49
20:QT:79:ARG:O	20:QT:83:ARG:HG3	2.12	0.49
25:RA:2271:G:OP1	46:R0:18:ALA:HB1	2.13	0.49
25:RA:2867:G:OP2	39:RT:119:LYS:NZ	2.34	0.49
31:RH:42:ARG:O	31:RH:52:VAL:HA	2.12	0.49
32:RI:74:ASN:OD1	32:RI:74:ASN:N	2.46	0.49
36:RQ:31:ASP:O	36:RQ:32:TYR:CG	2.66	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:RT:16:ARG:HD3	39:RT:19:LEU:HD11	1.94	0.49
46:R0:32:ARG:H	46:R0:35:ASN:ND2	2.11	0.49
51:R5:45:VAL:O	51:R5:45:VAL:HG12	2.13	0.49
51:R5:52:TYR:O	51:R5:53:ALA:CB	2.60	0.49
52:R6:25:LYS:HE2	52:R6:27:LYS:HD3	1.94	0.49
1:XA:64:G:H4'	1:XA:65:U:H5'	1.94	0.49
1:XA:263:A:OP2	20:XT:79:ARG:NH1	2.44	0.49
1:XA:431:A:H2'	1:XA:432:A:O4'	2.13	0.49
1:XA:737:A:H2'	1:XA:738:C:C6	2.48	0.49
1:XA:923:A:OP1	5:XE:21:ALA:HB2	2.13	0.49
5:XE:98:THR:HG22	5:XE:101:ILE:HD11	1.93	0.49
25:YA:27:G:O2'	25:YA:28:A:C8	2.66	0.49
25:YA:580:C:H2'	25:YA:581:C:C6	2.47	0.49
25:YA:782:A:N3	27:YD:226:MET:HB3	2.27	0.49
25:YA:2545:G:H2'	25:YA:2546:U:O4'	2.13	0.49
25:YA:2636:U:OP1	28:YE:79:ARG:HG3	2.12	0.49
27:YD:25:THR:O	27:YD:27:THR:HG22	2.12	0.49
35:YP:27:HIS:ND1	35:YP:27:HIS:N	2.61	0.49
38:YS:48:LEU:N	38:YS:48:LEU:CD1	2.76	0.49
54:Y8:33:ASN:O	54:Y8:35:GLN:N	2.46	0.49
54:Y8:35:GLN:OE1	54:Y8:35:GLN:HA	2.12	0.49
54:Y8:52:LYS:H	54:Y8:53:PRO:HD2	1.66	0.49
1:QA:1384:C:H2'	1:QA:1385:G:C8	2.48	0.48
2:QB:80:ILE:HG21	2:QB:212:GLN:HA	1.95	0.48
15:QO:87:ILE:HG22	15:QO:88:ARG:H	1.78	0.48
20:QT:33:ILE:HD12	20:QT:63:ILE:HG13	1.94	0.48
25:RA:1111:A:H4'	31:RH:3:ARG:CD	2.43	0.48
25:RA:1204:A:H2	25:RA:1241:A:N1	2.11	0.48
25:RA:1516:U:H2'	25:RA:1517:G:H8	1.77	0.48
25:RA:2528:U:OP1	55:R9:30:PRO:HG2	2.13	0.48
28:RE:38:THR:O	28:RE:42:ASP:HB2	2.13	0.48
28:RE:179:GLU:O	28:RE:180:ASN:HB2	2.12	0.48
31:RH:123:PHE:O	31:RH:125:VAL:HG23	2.13	0.48
31:RH:137:ASP:CB	31:RH:140:LYS:HB2	2.43	0.48
51:R5:56:LYS:HD2	51:R5:56:LYS:N	2.13	0.48
54:R8:29:LYS:O	54:R8:30:ARG:C	2.56	0.48
1:XA:545:C:OP1	4:XD:61:LYS:NZ	2.45	0.48
11:XK:59:TYR:CZ	11:XK:63:LEU:HD11	2.48	0.48
16:XP:26:ARG:HH21	16:XP:31:LYS:HB3	1.76	0.48
25:YA:530:G:O2'	25:YA:2021:C:O2'	2.30	0.48
25:YA:2392:A:H1'	35:YP:60:MET:HE3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2540:C:H2'	25:YA:2541:A:O4'	2.12	0.48
25:YA:2567:G:H2'	25:YA:2568:C:C6	2.48	0.48
35:YP:59:LEU:CA	35:YP:61:ARG:HE	2.24	0.48
35:YP:62:LEU:HD21	54:Y8:25:MET:CB	2.39	0.48
36:YQ:23:GLY:O	36:YQ:24:GLY:O	2.30	0.48
38:YS:55:ALA:O	38:YS:56:LEU:HB3	2.14	0.48
42:YW:51:LEU:HD23	42:YW:105:VAL:HG11	1.94	0.48
48:Y2:18:PRO:C	48:Y2:20:GLU:N	2.70	0.48
55:Y9:1:MET:HE1	55:Y9:32:HIS:CE1	2.48	0.48
1:QA:890:G:O2'	1:QA:906:G:O6	2.29	0.48
1:QA:1285:A:H5'	1:QA:1286:A:C2	2.47	0.48
15:QO:16:ALA:HB1	15:QO:21:ASP:HB3	1.95	0.48
21:QU:5:ASP:O	21:QU:11:GLY:HA3	2.13	0.48
25:RA:2777:G:OP2	25:RA:2781:A:O2'	2.20	0.48
31:RH:109:PHE:C	31:RH:111:HIS:H	2.21	0.48
43:RX:39:ILE:O	43:RX:43:VAL:HG12	2.13	0.48
44:RY:47:LYS:HG2	44:RY:60:PHE:CD1	2.48	0.48
51:R5:3:LYS:O	51:R5:4:HIS:C	2.56	0.48
1:XA:312:C:H2'	1:XA:313:A:C8	2.48	0.48
1:XA:957:U:N3	1:XA:960:U:OP2	2.32	0.48
1:XA:1308:U:OP1	13:XM:98:VAL:HG23	2.12	0.48
3:XC:130:VAL:HG21	3:XC:157:ILE:HG23	1.94	0.48
25:YA:2271:G:OP1	46:Y0:18:ALA:HB1	2.13	0.48
26:YB:40:U:H3	26:YB:43:C:H5"	1.77	0.48
30:YG:166:ASP:HA	30:YG:169:ALA:HB3	1.94	0.48
31:YH:13:LYS:HE2	31:YH:13:LYS:CA	2.40	0.48
31:YH:42:ARG:O	31:YH:52:VAL:HA	2.13	0.48
31:YH:137:ASP:CB	31:YH:140:LYS:HB2	2.43	0.48
34:YO:76:ALA:HB3	39:YT:75:ILE:HD12	1.95	0.48
36:YQ:90:VAL:C	36:YQ:92:GLY:N	2.71	0.48
40:YU:98:LEU:O	40:YU:102:GLU:N	2.37	0.48
45:YZ:5:LEU:HD11	45:YZ:39:VAL:HB	1.95	0.48
45:YZ:69:THR:HG22	45:YZ:90:VAL:HG22	1.96	0.48
1:QA:34:C:H2'	1:QA:35:G:C8	2.49	0.48
1:QA:148:G:H1	1:QA:174:C:N4	2.09	0.48
1:QA:266:G:H5"	1:QA:267:C:C5	2.48	0.48
1:QA:1128:C:H1'	1:QA:1146:A:H61	1.77	0.48
1:QA:1152:A:OP1	10:QJ:68:HIS:NE2	2.46	0.48
1:QA:1275:A:H2'	1:QA:1276:G:O4'	2.13	0.48
3:QC:162:GLN:NE2	3:QC:162:GLN:CA	2.68	0.48
7:QG:113:GLU:HG3	7:QG:119:ARG:HG2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:QL:38:THR:CG2	12:QL:57:LYS:HB3	2.44	0.48
19:QS:77:THR:HG22	19:QS:78:ARG:HD3	1.96	0.48
25:RA:1332:G:N2	25:RA:1609:A:O2'	2.46	0.48
25:RA:1783:A:HO2'	25:RA:2607:G:HO2'	1.60	0.48
25:RA:2292:C:P	38:RS:17:ARG:HH22	2.35	0.48
28:RE:77:ILE:CD1	28:RE:78:LEU:N	2.70	0.48
36:RQ:34:LEU:HD23	36:RQ:104:PHE:HD2	1.77	0.48
37:RR:97:VAL:HG22	37:RR:114:VAL:CG2	2.43	0.48
50:R4:42:PHE:CD1	50:R4:42:PHE:C	2.90	0.48
54:R8:41:ILE:HG13	54:R8:42:ARG:N	2.28	0.48
1:XA:953:G:H2'	1:XA:954:G:O4'	2.13	0.48
1:XA:1060:C:C5	3:XC:2:GLY:HA2	2.48	0.48
3:XC:79:ARG:HH12	3:XC:82:GLU:HG3	1.78	0.48
3:XC:134:ILE:HG23	3:XC:151:VAL:HB	1.94	0.48
5:XE:50:GLU:HG3	5:XE:52:PRO:HD2	1.95	0.48
6:XF:35:ALA:HA	6:XF:67:MET:HB3	1.94	0.48
12:XL:85:ILE:HD11	12:XL:98:TYR:CB	2.42	0.48
15:XO:70:LEU:HD11	15:XO:77:ARG:HG3	1.96	0.48
22:XV:14:G:N2	22:XV:48:C:H42	2.09	0.48
25:YA:305:U:H2'	25:YA:306:U:C6	2.48	0.48
25:YA:1062:G:H2'	25:YA:1063:G:C8	2.47	0.48
25:YA:2212:A:H1'	25:YA:2215:G:C5	2.48	0.48
25:YA:2401:U:H2'	25:YA:2402:C:H5"	1.95	0.48
26:YB:74:U:H2'	26:YB:75:G:O4'	2.13	0.48
27:YD:48:ARG:HG3	27:YD:48:ARG:HH11	1.78	0.48
28:YE:38:THR:O	28:YE:42:ASP:HB2	2.13	0.48
31:YH:53:GLU:CD	31:YH:54:ARG:H	2.21	0.48
31:YH:109:PHE:C	31:YH:111:HIS:H	2.21	0.48
38:YS:33:LYS:HB3	38:YS:34:HIS:CD2	2.48	0.48
1:QA:110:C:O2'	16:QP:25:ARG:O	2.26	0.48
1:QA:1218:C:H2'	1:QA:1219:U:C6	2.48	0.48
2:QB:163:PHE:HD1	2:QB:185:ILE:HG13	1.78	0.48
10:QJ:32:ALA:HB3	10:QJ:76:ASN:HB2	1.96	0.48
13:QM:23:TYR:HE1	13:QM:70:LEU:HD12	1.77	0.48
19:QS:26:GLY:O	19:QS:28:LYS:N	2.41	0.48
25:RA:18:C:O2'	25:RA:553:U:OP1	2.30	0.48
25:RA:27:G:O2'	25:RA:28:A:H8	1.96	0.48
25:RA:443:A:N7	29:RF:45:ARG:HD2	2.27	0.48
25:RA:1045:A:N9	25:RA:1111:A:N6	2.62	0.48
25:RA:1047:G:H2'	25:RA:1110:G:H1	1.77	0.48
25:RA:1436:G:H1'	25:RA:1477:A:O2'	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2096:U:H3	25:RA:2193:G:H1	1.61	0.48
25:RA:2250:G:C5	36:RQ:82:ARG:HD2	2.48	0.48
28:RE:78:LEU:CD2	28:RE:79:ARG:HD2	2.44	0.48
30:RG:98:ARG:HE	30:RG:98:ARG:HB2	1.38	0.48
31:RH:104:GLU:HG3	31:RH:114:VAL:HG22	1.96	0.48
31:RH:119:GLU:CD	31:RH:120:GLY:H	2.22	0.48
32:RI:53:ALA:O	32:RI:57:ARG:HG3	2.08	0.48
33:RN:4:TYR:O	40:RU:64:ARG:NH1	2.46	0.48
47:R1:53:VAL:HB	47:R1:58:ILE:HD12	1.94	0.48
49:R3:4:LEU:O	49:R3:36:VAL:HA	2.13	0.48
54:R8:56:GLU:O	54:R8:58:ILE:N	2.47	0.48
1:XA:429:U:H1'	1:XA:430:A:H5"	1.94	0.48
1:XA:686:U:O2'	11:XK:42:TRP:NE1	2.46	0.48
2:XB:101:MET:HA	2:XB:108:ILE:HG13	1.95	0.48
3:XC:7:PRO:O	3:XC:11:ARG:HG2	2.13	0.48
6:XF:36:ARG:CZ	6:XF:38:GLU:HG2	2.44	0.48
12:XL:38:THR:CG2	12:XL:57:LYS:HB3	2.44	0.48
25:YA:242:G:C3'	54:Y8:6:THR:HG23	2.43	0.48
25:YA:910:A:N3	25:YA:2264:C:O2'	2.43	0.48
25:YA:1590:U:H2'	25:YA:1591:G:C8	2.49	0.48
25:YA:2439:A:H5'	25:YA:2439:A:H8	1.78	0.48
27:YD:45:ASN:CG	27:YD:46:GLN:N	2.68	0.48
29:YF:155:LEU:HD23	29:YF:186:ILE:HA	1.95	0.48
31:YH:82:GLY:O	31:YH:83:TYR:O	2.31	0.48
47:Y1:41:ARG:HH11	47:Y1:41:ARG:HG3	1.79	0.48
1:QA:411:A:C5	1:QA:413:G:H1'	2.48	0.48
1:QA:630:G:N2	1:QA:631:G:H21	2.10	0.48
1:QA:718:G:H5'	11:QK:117:ASN:ND2	2.28	0.48
3:QC:150:LYS:HG3	3:QC:169:ALA:HB2	1.96	0.48
4:QD:166:LYS:HD3	27:YD:134:ARG:HH12	1.78	0.48
10:QJ:80:LYS:HA	10:QJ:80:LYS:HD3	1.69	0.48
13:QM:49:THR:HG22	13:QM:51:ALA:H	1.79	0.48
25:RA:2298:A:H2'	25:RA:2299:G:O4'	2.13	0.48
25:RA:2495:G:H5"	36:RQ:81:VAL:CG1	2.44	0.48
36:RQ:21:THR:HB	36:RQ:22:LYS:H	1.42	0.48
44:RY:81:LYS:NZ	44:RY:98:VAL:HG11	2.28	0.48
45:RZ:45:ASP:O	45:RZ:49:ARG:HG2	2.13	0.48
1:XA:1175:G:H2'	1:XA:1176:A:C8	2.48	0.48
2:XB:37:ASN:C	2:XB:39:ILE:H	2.20	0.48
11:XK:18:ARG:NH2	11:XK:35:PRO:O	2.45	0.48
25:YA:728:G:H4'	27:YD:13:ARG:HD2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1826:G:O2'	27:YD:242:ARG:NH2	2.46	0.48
25:YA:2467:C:C4'	36:YQ:123:HIS:CD2	2.96	0.48
25:YA:2756:U:OP2	55:Y9:19:ARG:NH2	2.46	0.48
27:YD:10:THR:O	27:YD:11:PRO:C	2.56	0.48
27:YD:33:LEU:HB3	27:YD:34:VAL:H	1.48	0.48
27:YD:44:ASN:H	27:YD:44:ASN:ND2	1.97	0.48
28:YE:69:LYS:C	28:YE:71:GLY:N	2.71	0.48
29:YF:107:LYS:O	29:YF:110:LEU:N	2.47	0.48
29:YF:128:ALA:O	29:YF:129:PHE:HB2	2.13	0.48
29:YF:198:ALA:O	29:YF:201:VAL:HG12	2.13	0.48
31:YH:7:LEU:N	31:YH:8:PRO:CD	2.77	0.48
31:YH:12:PRO:HD3	31:YH:48:GLY:O	2.13	0.48
31:YH:23:ARG:HD2	31:YH:34:GLU:OE2	2.12	0.48
43:YX:63:LYS:O	43:YX:64:LYS:HD2	2.14	0.48
51:Y5:41:PRO:O	51:Y5:44:THR:OG1	2.32	0.48
1:QA:1004:A:O5'	1:QA:1025:U:N3	2.46	0.48
7:QG:57:GLU:OE1	7:QG:57:GLU:N	2.41	0.48
10:QJ:50:ILE:HD12	10:QJ:57:LYS:HA	1.95	0.48
16:QP:43:LYS:HA	16:QP:48:TRP:HB3	1.96	0.48
25:RA:414:C:H2'	25:RA:415:A:H8	1.78	0.48
25:RA:1062:G:H2'	25:RA:1063:G:C8	2.48	0.48
25:RA:1286:A:O2'	25:RA:1288:U:OP2	2.25	0.48
29:RF:102:PRO:HB2	29:RF:105:VAL:HG23	1.94	0.48
35:RP:65:ARG:HH21	54:R8:46:ARG:HH12	1.60	0.48
37:RR:2:ARG:HA	37:RR:5:LYS:HE3	1.95	0.48
38:RS:64:GLU:O	38:RS:68:GLN:HG3	2.14	0.48
50:R4:48:ARG:C	50:R4:49:PHE:HD1	2.22	0.48
51:R5:49:CYS:SG	51:R5:58:LEU:HB2	2.53	0.48
54:R8:35:GLN:OE1	54:R8:35:GLN:HA	2.12	0.48
6:XF:10:LEU:HD22	6:XF:61:LEU:HD11	1.94	0.48
9:XI:9:ARG:HB2	9:XI:14:VAL:HA	1.96	0.48
12:XL:8:ASN:OD1	17:XQ:34:LYS:NZ	2.32	0.48
25:YA:1535:U:N3	25:YA:1537:C:H1'	2.27	0.48
27:YD:35:LYS:HD2	27:YD:104:TYR:CE1	2.49	0.48
28:YE:174:ASP:O	28:YE:182:LEU:HD12	2.14	0.48
31:YH:124:GLU:HB3	31:YH:132:ARG:CD	2.44	0.48
32:YI:133:HIS:HB2	32:YI:134:PRO:CD	2.43	0.48
36:YQ:31:ASP:O	36:YQ:32:TYR:CG	2.66	0.48
39:YT:102:ILE:HB	39:YT:110:ILE:HD13	1.95	0.48
41:YV:15:GLU:O	41:YV:18:LEU:HB2	2.14	0.48
44:YY:35:TYR:CD2	44:YY:69:ALA:HB3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:Y1:25:LYS:C	47:Y1:27:GLU:H	2.22	0.48
48:Y2:16:LEU:O	48:Y2:17:SER:CB	2.56	0.48
1:QA:620:C:C2	4:QD:135:LEU:HG	2.48	0.48
1:QA:857:C:H2'	1:QA:858:G:O4'	2.13	0.48
1:QA:1080:A:C5'	5:QE:16:THR:HG21	2.44	0.48
1:QA:1126:U:H1'	1:QA:1280:A:N7	2.29	0.48
1:QA:1343:G:H2'	1:QA:1344:C:C6	2.49	0.48
1:QA:1466:C:H2'	1:QA:1467:G:O4'	2.14	0.48
9:QI:17:VAL:HG11	9:QI:81:ILE:HA	1.95	0.48
11:QK:48:ILE:HD11	11:QK:64:ALA:HA	1.95	0.48
15:QO:26:GLU:H	15:QO:26:GLU:HG2	1.42	0.48
17:QQ:100:LYS:O	17:QQ:101:ARG:NE	2.47	0.48
20:QT:30:LYS:NZ	20:QT:80:ARG:NH1	2.62	0.48
25:RA:49:A:H4'	25:RA:50:U:H5"	1.94	0.48
25:RA:372:G:O2'	25:RA:373:U:P	2.71	0.48
25:RA:492:A:H2'	25:RA:493:G:O4'	2.13	0.48
25:RA:780:G:H21	25:RA:783:A:H62	1.60	0.48
25:RA:1782:C:H1'	25:RA:2609:U:H5"	1.95	0.48
28:RE:61:ARG:CB	28:RE:62:PRO:CD	2.90	0.48
29:RF:103:LYS:HA	29:RF:106:ARG:HG3	1.94	0.48
30:RG:81:LYS:O	30:RG:82:LEU:HB2	2.13	0.48
30:RG:95:ARG:C	30:RG:99:MET:HG2	2.38	0.48
31:RH:7:LEU:N	31:RH:8:PRO:CD	2.77	0.48
44:RY:51:VAL:O	44:RY:56:PRO:HA	2.14	0.48
50:R4:60:GLN:O	50:R4:63:TYR:HB3	2.14	0.48
2:XB:74:LYS:HE2	2:XB:169:LYS:HG3	1.94	0.48
25:YA:127:A:H5"	25:YA:128:C:C6	2.48	0.48
25:YA:1678:G:H8	25:YA:1678:G:O5'	1.96	0.48
25:YA:1889:A:O2'	25:YA:2087:G:H5'	2.13	0.48
25:YA:1930:G:N2	25:YA:1969:A:OP2	2.36	0.48
25:YA:2023:G:H4'	25:YA:2617:C:O3'	2.14	0.48
25:YA:2563:U:H1'	25:YA:2566:A:N6	2.28	0.48
25:YA:2811:G:OP1	28:YE:61:ARG:CG	2.62	0.48
27:YD:12:SER:C	27:YD:14:ARG:N	2.70	0.48
27:YD:130:ALA:HA	27:YD:192:THR:HA	1.95	0.48
28:YE:77:ILE:CD1	28:YE:78:LEU:N	2.70	0.48
31:YH:41:MET:HG3	31:YH:54:ARG:HA	1.96	0.48
31:YH:127:GLU:HB3	31:YH:128:PRO:HD2	1.92	0.48
31:YH:153:LYS:CB	31:YH:154:PRO:CD	2.69	0.48
36:YQ:42:ILE:HD12	36:YQ:42:ILE:N	2.29	0.48
38:YS:18:ILE:O	38:YS:19:LYS:O	2.30	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:YU:60:LEU:HD22	40:YU:60:LEU:O	2.14	0.48
40:YU:97:ASP:OD2	40:YU:101:ARG:NH1	2.46	0.48
48:Y2:33:MET:O	48:Y2:37:PHE:HD1	1.95	0.48
48:Y2:69:ARG:HH11	48:Y2:69:ARG:HB3	1.79	0.48
1:QA:35:G:H2'	1:QA:36:C:C6	2.49	0.48
1:QA:164:U:H2'	1:QA:165:C:C6	2.49	0.48
9:QI:40:LEU:O	9:QI:42:ARG:N	2.46	0.48
12:QL:22:SER:C	12:QL:24:VAL:H	2.22	0.48
13:QM:121:LYS:NZ	24:QY:40:G:H4'	2.29	0.48
20:QT:12:ALA:O	20:QT:15:ARG:HB2	2.14	0.48
25:RA:245:G:O2'	25:RA:384:U:O2	2.28	0.48
25:RA:263:C:H2'	25:RA:264:C:O4'	2.14	0.48
25:RA:1534:G:N3	25:RA:1534:G:H2'	2.27	0.48
25:RA:2283:C:P	52:R6:5:VAL:HG13	2.54	0.48
25:RA:2376:A:H2'	25:RA:2377:A:O4'	2.12	0.48
30:RG:82:LEU:HD21	30:RG:88:ILE:HG13	1.96	0.48
32:RI:29:TYR:HD2	32:RI:30:LEU:HD23	1.79	0.48
36:RQ:87:LYS:O	36:RQ:89:ASN:N	2.43	0.48
45:RZ:111:VAL:HG22	45:RZ:112:ARG:N	2.19	0.48
54:R8:30:ARG:O	54:R8:31:HIS:HB2	2.12	0.48
54:R8:40:GLU:O	54:R8:41:ILE:C	2.56	0.48
1:XA:1318:A:H5'	19:XS:11:VAL:HG11	1.96	0.48
9:XI:114:TYR:N	9:XI:114:TYR:CD1	2.81	0.48
25:YA:330:A:HO2'	25:YA:331:A:H8	1.58	0.48
25:YA:586:A:H5'	29:YF:89:VAL:HG21	1.96	0.48
25:YA:774:A:H2	25:YA:787:U:O2'	1.97	0.48
25:YA:1952:A:C6	25:YA:1953:A:N1	2.81	0.48
25:YA:2059:A:H5'	25:YA:2060:A:OP2	2.14	0.48
25:YA:2119:A:H61	25:YA:2168:G:H22	1.60	0.48
25:YA:2349:G:OP2	54:Y8:42:ARG:HD3	2.14	0.48
25:YA:2619:C:O2'	28:YE:156:MET:CE	2.62	0.48
27:YD:35:LYS:CG	27:YD:64:ILE:CG2	2.92	0.48
28:YE:23:VAL:HG12	28:YE:173:VAL:HG21	1.94	0.48
29:YF:45:ARG:NH1	29:YF:45:ARG:HG2	2.28	0.48
29:YF:129:PHE:CD2	29:YF:163:VAL:HG21	2.48	0.48
31:YH:119:GLU:CD	31:YH:120:GLY:H	2.22	0.48
31:YH:120:GLY:HA3	31:YH:140:LYS:NZ	2.27	0.48
36:YQ:19:GLY:O	36:YQ:98:LYS:HD3	2.14	0.48
38:YS:66:ALA:HA	38:YS:69:VAL:HG12	1.96	0.48
41:YV:76:LYS:HB2	41:YV:81:TYR:HB3	1.95	0.48
42:YW:67:ASP:N	42:YW:67:ASP:OD1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:Y2:27:GLU:H	48:Y2:27:GLU:CD	2.17	0.48
50:Y4:43:TYR:CD1	50:Y4:43:TYR:C	2.92	0.48
52:Y6:27:LYS:NZ	52:Y6:27:LYS:HB2	2.28	0.48
1:QA:42:G:H2'	1:QA:43:C:O4'	2.13	0.48
1:QA:520:A:OP2	12:QL:51:ALA:HB1	2.13	0.48
25:RA:2404:C:H1'	35:RP:67:MET:HE1	1.94	0.48
28:RE:93:VAL:H	28:RE:95:ILE:CD1	2.23	0.48
28:RE:203:LYS:HD2	28:RE:203:LYS:C	2.38	0.48
29:RF:133:ASN:O	29:RF:135:LYS:N	2.46	0.48
31:RH:131:VAL:HG12	31:RH:132:ARG:N	2.29	0.48
36:RQ:19:GLY:O	36:RQ:98:LYS:HD3	2.14	0.48
36:RQ:119:ARG:O	36:RQ:123:HIS:HD2	1.97	0.48
52:R6:18:ARG:HB2	52:R6:44:ARG:HH12	1.77	0.48
1:XA:263:A:P	20:XT:79:ARG:HH11	2.37	0.48
1:XA:1117:G:H5"	9:XI:104:ARG:NH1	2.29	0.48
5:XE:41:VAL:HG13	5:XE:113:ALA:CB	2.36	0.48
12:XL:22:SER:C	12:XL:24:VAL:H	2.21	0.48
12:XL:27:LEU:C	12:XL:29:GLY:H	2.20	0.48
19:XS:39:THR:HG22	19:XS:40:ILE:H	1.78	0.48
20:XT:93:GLU:OE1	20:XT:94:ALA:N	2.46	0.48
20:XT:98:PRO:C	20:XT:100:ILE:H	2.21	0.48
25:YA:226:G:H2'	25:YA:227:A:C8	2.49	0.48
25:YA:1007:C:OP1	33:YN:35:ARG:NH1	2.46	0.48
25:YA:1022:G:OP2	33:YN:65:LYS:HE3	2.14	0.48
25:YA:1533:C:N4	25:YA:1538:G:H1	2.11	0.48
28:YE:15:PHE:CD1	28:YE:20:ALA:HB2	2.49	0.48
28:YE:55:ASN:O	28:YE:57:LYS:N	2.44	0.48
28:YE:119:ARG:HD3	28:YE:160:TYR:HD2	1.78	0.48
1:QA:299:G:H2'	1:QA:300:A:C8	2.49	0.48
1:QA:381:C:H2'	1:QA:382:A:O4'	2.14	0.48
1:QA:793:U:O2	1:QA:1516:G:H4'	2.13	0.48
1:QA:1143:G:H2'	1:QA:1144:G:H8	1.79	0.48
1:QA:1499:A:H1'	1:QA:1520:G:H5'	1.94	0.48
8:QH:82:HIS:HD1	8:QH:138:TRP:NE1	2.11	0.48
8:QH:102:ARG:NH1	8:QH:105:ARG:HH22	2.12	0.48
10:QJ:78:ASN:O	10:QJ:82:ILE:HG12	2.14	0.48
25:RA:526:A:OP1	25:RA:527:C:OP1	2.31	0.48
25:RA:747:U:C2	51:R5:2:ALA:CB	2.96	0.48
25:RA:817:C:O2'	25:RA:839:U:H5"	2.13	0.48
25:RA:1678:G:H22	25:RA:1989:G:N2	2.12	0.48
25:RA:2126:A:N6	25:RA:2163:C:O2'	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:RD:25:THR:O	27:RD:27:THR:HG22	2.14	0.48
35:RP:61:ARG:HG3	54:R8:13:ARG:HH11	1.79	0.48
43:RX:83:VAL:HG11	43:RX:87:GLN:HB2	1.96	0.48
1:XA:522:C:H41	12:XL:53:ARG:HH22	1.61	0.48
1:XA:523:A:N6	12:XL:92:ASP:HB2	2.27	0.48
1:XA:790:A:C6	1:XA:791:G:C6	3.02	0.48
1:XA:938:A:N3	1:XA:1376:U:O2'	2.46	0.48
1:XA:1489:G:H2'	1:XA:1490:C:O4'	2.14	0.48
9:XI:91:ASP:C	9:XI:93:ARG:H	2.21	0.48
16:XP:22:THR:HA	16:XP:33:ILE:HG12	1.96	0.48
24:XY:37:A:O2'	25:YA:1913:A:N1	2.45	0.48
25:YA:592:G:C2	54:Y8:4:MET:HE2	2.49	0.48
25:YA:919:G:N2	25:YA:2269:A:OP2	2.46	0.48
25:YA:1178:C:H2'	25:YA:1179:C:C6	2.49	0.48
25:YA:1654:A:OP2	37:YR:2:ARG:HD2	2.14	0.48
25:YA:2126:A:N6	25:YA:2163:C:O2'	2.47	0.48
25:YA:2397:G:H5"	47:Y1:28:GLY:HA2	1.96	0.48
26:YB:104:A:H5'	45:YZ:72:ARG:HD3	1.96	0.48
27:YD:27:THR:O	27:YD:29:PRO:CD	2.62	0.48
27:YD:198:ASN:C	27:YD:198:ASN:ND2	2.69	0.48
36:YQ:60:ARG:NH2	36:YQ:60:ARG:HB2	2.28	0.48
1:QA:335:C:O2'	1:QA:1433:A:N3	2.45	0.47
2:QB:97:TRP:CH2	2:QB:173:ALA:HA	2.49	0.47
12:QL:27:LEU:C	12:QL:29:GLY:H	2.20	0.47
12:QL:120:TYR:O	12:QL:121:GLY:C	2.57	0.47
25:RA:78:A:H2'	25:RA:79:G:H8	1.79	0.47
25:RA:593:G:H1	25:RA:664:C:N4	2.10	0.47
25:RA:922:U:H2'	25:RA:923:C:C6	2.49	0.47
25:RA:1790:C:H5"	25:RA:1791:A:OP1	2.14	0.47
28:RE:69:LYS:C	28:RE:71:GLY:N	2.71	0.47
33:RN:7:LYS:H	33:RN:7:LYS:HD2	1.79	0.47
36:RQ:42:ILE:HD12	36:RQ:42:ILE:N	2.29	0.47
1:XA:1227:A:OP1	19:XS:80:TYR:OH	2.25	0.47
1:XA:1312:G:H5"	50:Y4:67:TYR:OH	2.14	0.47
3:XC:34:LEU:HD23	3:XC:38:ARG:HG3	1.95	0.47
5:XE:6:PHE:CE1	5:XE:36:ASP:HB3	2.48	0.47
8:XH:49:GLU:HG2	8:XH:62:TYR:HE2	1.78	0.47
10:XJ:38:ILE:HD11	10:XJ:71:LEU:HD23	1.96	0.47
12:XL:52:LEU:O	12:XL:54:LYS:NZ	2.30	0.47
25:YA:273:G:H1	25:YA:364:C:N4	2.11	0.47
25:YA:286:C:H2'	25:YA:287:C:C6	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1614:A:H61	42:YW:88:ARG:H	1.62	0.47
31:YH:45:VAL:O	31:YH:45:VAL:HG13	2.14	0.47
33:YN:134:ARG:N	33:YN:135:PRO:HD3	2.29	0.47
41:YV:44:LYS:O	41:YV:46:VAL:HG12	2.13	0.47
44:YY:44:ILE:HG13	44:YY:45:VAL:N	2.28	0.47
45:YZ:10:ARG:HH21	45:YZ:26:GLY:H	1.61	0.47
45:YZ:182:LYS:CG	45:YZ:183:LEU:HD23	2.39	0.47
51:Y5:46:CYS:O	51:Y5:48:GLU:N	2.38	0.47
54:Y8:53:PRO:CD	54:Y8:54:GLU:N	2.77	0.47
1:QA:617:G:H1	1:QA:623:C:N4	2.09	0.47
1:QA:1052:U:H5"	1:QA:1053:G:OP2	2.14	0.47
1:QA:1225:A:H5"	1:QA:1226:C:OP2	2.13	0.47
1:QA:1305:G:O2'	1:QA:1306:A:H8	1.97	0.47
1:QA:1305:G:O2'	1:QA:1306:A:O4'	2.32	0.47
2:QB:25:ASN:O	2:QB:27:LYS:N	2.47	0.47
6:QF:41:GLU:HB2	6:QF:62:TRP:CE3	2.50	0.47
12:QL:85:ILE:HD11	12:QL:98:TYR:CB	2.43	0.47
12:QL:127:GLU:O	12:QL:128:ALA:CB	2.62	0.47
20:QT:30:LYS:O	20:QT:33:ILE:HB	2.14	0.47
25:RA:607:U:H5	25:RA:619:G:C5	2.31	0.47
25:RA:781:A:H2	25:RA:1776:G:N3	2.12	0.47
25:RA:1041:C:H42	25:RA:1114:G:H1	1.61	0.47
28:RE:18:ASP:O	28:RE:19:ARG:C	2.56	0.47
28:RE:120:TRP:O	28:RE:121:ASN:HB2	2.14	0.47
31:RH:45:VAL:HG13	31:RH:45:VAL:O	2.14	0.47
31:RH:154:PRO:CG	31:RH:162:ILE:O	2.61	0.47
32:RI:31:LEU:HD11	32:RI:38:LEU:HG	1.96	0.47
36:RQ:57:HIS:ND1	36:RQ:58:PHE:N	2.62	0.47
38:RS:88:ASP:CG	38:RS:89:ARG:H	2.21	0.47
55:R9:27:CYS:SG	55:R9:32:HIS:HB2	2.55	0.47
1:XA:164:U:H2'	1:XA:165:C:C6	2.49	0.47
1:XA:166:G:H2'	1:XA:167:G:H8	1.79	0.47
1:XA:228:A:H2'	1:XA:229:U:O4'	2.13	0.47
1:XA:406:G:H5'	4:XD:5:ILE:HD13	1.96	0.47
3:XC:70:VAL:HG12	3:XC:72:LYS:H	1.79	0.47
5:XE:110:LEU:HD13	5:XE:118:ILE:HG21	1.95	0.47
5:XE:152:ARG:HG2	8:XH:79:VAL:HG13	1.96	0.47
9:XI:121:ARG:NH1	9:XI:122:ALA:O	2.47	0.47
10:XJ:55:LYS:HG3	10:XJ:56:HIS:H	1.78	0.47
18:XR:36:ASN:O	18:XR:36:ASN:ND2	2.41	0.47
25:YA:1427:A:H4'	25:YA:1428:C:O5'	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2277:G:H5”	36:YQ:85:LYS:HG3	1.96	0.47
28:YE:61:ARG:CB	28:YE:62:PRO:HD3	2.41	0.47
29:YF:155:LEU:HA	29:YF:174:VAL:HG12	1.95	0.47
45:YZ:15:PRO:O	45:YZ:19:ARG:HB2	2.13	0.47
1:QA:45:U:H2’	1:QA:46:G:C8	2.49	0.47
1:QA:1039:C:H2’	1:QA:1040:U:O4’	2.14	0.47
4:QD:107:ARG:HH21	4:QD:194:LEU:HD21	1.79	0.47
25:RA:498:G:N3	44:RY:47:LYS:NZ	2.62	0.47
25:RA:871:U:H4’	36:RQ:69:PHE:CE2	2.49	0.47
25:RA:1090:U:C2	25:RA:1102:C:H1’	2.49	0.47
25:RA:1289:C:H2’	25:RA:1290:C:C6	2.49	0.47
25:RA:1417:C:H2’	25:RA:1418:G:O4’	2.14	0.47
25:RA:1817:G:OP1	27:RD:88:ARG:NH2	2.45	0.47
25:RA:2298:A:C2	25:RA:2299:G:H1’	2.50	0.47
27:RD:35:LYS:HE3	27:RD:64:ILE:C	2.39	0.47
28:RE:56:PRO:O	28:RE:57:LYS:CB	2.61	0.47
28:RE:174:ASP:O	28:RE:182:LEU:HD12	2.14	0.47
30:RG:5:VAL:H	50:R4:25:TYR:HE2	1.62	0.47
30:RG:145:THR:O	30:RG:147:ASP:N	2.47	0.47
31:RH:82:GLY:O	31:RH:83:TYR:O	2.31	0.47
31:RH:124:GLU:HB3	31:RH:132:ARG:CD	2.44	0.47
35:RP:61:ARG:NE	54:R8:13:ARG:CD	2.60	0.47
47:R1:91:LYS:O	47:R1:94:LEU:N	2.36	0.47
51:R5:57:VAL:O	51:R5:57:VAL:HG13	2.14	0.47
1:XA:1213:A:N6	1:XA:1215:G:N3	2.62	0.47
2:XB:84:GLU:OE1	2:XB:87:ARG:NH2	2.43	0.47
3:XC:81:GLY:O	3:XC:85:ARG:HB2	2.14	0.47
5:XE:11:ILE:HG21	5:XE:108:ALA:HB2	1.96	0.47
9:XI:126:SER:O	9:XI:128:ARG:N	2.43	0.47
12:XL:50:SER:O	12:XL:51:ALA:HB2	2.14	0.47
20:XT:35:THR:O	20:XT:39:LYS:HG3	2.14	0.47
25:YA:910:A:N7	36:YQ:13:GLN:HG3	2.29	0.47
31:YH:131:VAL:HG12	31:YH:132:ARG:N	2.29	0.47
38:YS:40:ILE:HG22	38:YS:41:ASP:N	2.28	0.47
38:YS:105:ALA:C	38:YS:110:LEU:HD21	2.38	0.47
39:YT:11:GLU:OE1	39:YT:11:GLU:N	2.43	0.47
47:Y1:91:LYS:HB3	47:Y1:92:LYS:H	1.44	0.47
54:Y8:41:ILE:HG13	54:Y8:42:ARG:N	2.28	0.47
54:Y8:56:GLU:O	54:Y8:58:ILE:N	2.47	0.47
1:QA:429:U:H3’	4:QD:22:LYS:HZ1	1.79	0.47
2:QB:211:ILE:O	2:QB:215:LEU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:QE:101:ILE:N	5:QE:101:ILE:CD1	2.73	0.47
9:QI:28:VAL:HG22	9:QI:63:ILE:HB	1.95	0.47
12:QL:15:ARG:HG3	17:QQ:32:TYR:OH	2.14	0.47
13:QM:57:ARG:HB2	13:QM:57:ARG:HH11	1.80	0.47
14:QN:41:ARG:CZ	14:QN:42:ILE:HD11	2.44	0.47
17:QQ:76:LEU:HD21	17:QQ:79:SER:HB2	1.97	0.47
25:RA:270(E):G:H1	25:RA:270(U):C:H42	1.63	0.47
25:RA:467:G:OP1	53:R7:33:ARG:NH1	2.45	0.47
25:RA:1654:A:H2	28:RE:113:PHE:CD2	2.33	0.47
25:RA:1803:A:N1	25:RA:1822:G:O2'	2.42	0.47
28:RE:55:ASN:C	28:RE:57:LYS:N	2.62	0.47
31:RH:67:LEU:O	31:RH:71:LEU:HB2	2.14	0.47
48:R2:41:ILE:HD11	48:R2:44:LEU:HD12	1.96	0.47
1:XA:762:C:H2'	1:XA:763:G:H8	1.78	0.47
1:XA:891:U:H2'	1:XA:892:A:H8	1.79	0.47
1:XA:1186:G:H21	14:XN:61:TRP:C	2.22	0.47
1:XA:1280:A:O2'	1:XA:1281:U:OP1	2.27	0.47
4:XD:15:GLU:HG2	4:XD:63:LYS:HB2	1.97	0.47
12:XL:43:VAL:HG23	12:XL:93:LEU:HD22	1.97	0.47
22:XV:14:G:H22	22:XV:48:C:N4	2.11	0.47
25:YA:519:U:H2'	25:YA:520:G:C8	2.50	0.47
25:YA:1278:A:H4'	37:YR:34:ILE:HD12	1.95	0.47
25:YA:2420:C:N4	54:Y8:30:ARG:HD2	2.29	0.47
25:YA:2452:C:O4'	56:Z6:101:PPU:HB2	2.13	0.47
25:YA:2579:C:H2'	25:YA:2580:U:O4'	2.15	0.47
28:YE:65:GLY:HA2	28:YE:70:ALA:HB3	1.95	0.47
31:YH:127:GLU:OE2	31:YH:130:ARG:NH2	2.48	0.47
31:YH:154:PRO:CG	31:YH:162:ILE:O	2.61	0.47
35:YP:101:VAL:C	35:YP:103:ALA:H	2.23	0.47
42:YW:110:LYS:HG3	42:YW:111:HIS:H	1.80	0.47
1:QA:137:C:H42	1:QA:226:G:H1	1.62	0.47
1:QA:1244:C:H42	1:QA:1293:G:H1	1.61	0.47
2:QB:70:PHE:O	2:QB:93:VAL:N	2.48	0.47
6:QF:61:LEU:HB3	6:QF:63:TYR:HE1	1.80	0.47
7:QG:99:LEU:HD22	7:QG:103:TRP:CZ2	2.49	0.47
10:QJ:55:LYS:CG	10:QJ:56:HIS:H	2.27	0.47
12:QL:50:SER:O	12:QL:51:ALA:HB2	2.14	0.47
25:RA:530:G:O2'	25:RA:2021:C:O2'	2.33	0.47
25:RA:605:C:O2	25:RA:657:U:O2'	2.33	0.47
25:RA:1492:G:OP1	25:RA:2210:G:N1	2.43	0.47
25:RA:2816:C:O3'	37:RR:99:LYS:NZ	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:RE:195:LEU:HD12	28:RE:196:VAL:N	2.29	0.47
40:RU:97:ASP:OD2	40:RU:101:ARG:NH1	2.48	0.47
44:RY:21:LYS:HG3	44:RY:22:GLY:N	2.30	0.47
50:R4:8:LYS:O	50:R4:9:LEU:CB	2.62	0.47
1:XA:376:G:H5"	16:XP:5:ARG:HB2	1.97	0.47
1:XA:1000:A:H2'	1:XA:1001:G:H8	1.79	0.47
1:XA:1286:A:H2	21:XU:22:ARG:HH22	1.62	0.47
1:XA:1306:A:H61	1:XA:1331:G:H1'	1.80	0.47
1:XA:1306:A:N6	1:XA:1331:G:H1'	2.30	0.47
3:XC:79:ARG:NH1	3:XC:82:GLU:HG3	2.29	0.47
5:XE:8:GLU:OE2	5:XE:63:ARG:NH2	2.46	0.47
6:XF:69:GLU:O	6:XF:72:VAL:HG12	2.14	0.47
9:XI:4:TYR:CZ	9:XI:88:TYR:HB2	2.49	0.47
10:XJ:53:PRO:O	14:XN:41:ARG:NH2	2.39	0.47
18:XR:66:LEU:O	18:XR:70:ILE:HG13	2.15	0.47
25:YA:141(A):C:H2'	25:YA:142:G:O4'	2.14	0.47
25:YA:270(F):U:H2'	25:YA:270(G):C:C6	2.50	0.47
25:YA:1048:A:P	25:YA:1110:G:H22	2.37	0.47
25:YA:1496:A:H5'	25:YA:1497:U:OP1	2.14	0.47
25:YA:1655:A:H3'	25:YA:1656:C:C6	2.49	0.47
25:YA:2688:U:H1'	25:YA:2721:A:N6	2.30	0.47
27:YD:31:LYS:C	27:YD:32:SER:O	2.54	0.47
28:YE:56:PRO:O	28:YE:57:LYS:CB	2.61	0.47
35:YP:68:GLN:HG2	54:Y8:12:LYS:HG2	1.97	0.47
36:YQ:80:GLU:HG3	36:YQ:81:VAL:N	2.27	0.47
47:Y1:53:VAL:HG22	47:Y1:74:VAL:HG13	1.96	0.47
48:Y2:6:VAL:O	48:Y2:7:ARG:C	2.57	0.47
1:QA:281:G:H8	1:QA:281:G:OP2	1.97	0.47
1:QA:457:C:H2'	1:QA:458:C:H6	1.80	0.47
1:QA:524:G:H2'	1:QA:525:C:C6	2.49	0.47
1:QA:1129:C:H5'	1:QA:1130:A:H5'	1.95	0.47
2:QB:85:ALA:HB3	2:QB:92:TYR:HD2	1.80	0.47
13:QM:8:GLU:CD	30:RG:115:ARG:HD3	2.40	0.47
25:RA:262:A:H2'	25:RA:263:C:O4'	2.15	0.47
25:RA:859:G:O2'	25:RA:860:U:P	2.72	0.47
25:RA:1711:C:N4	25:RA:1747:G:H1	2.10	0.47
25:RA:2114:A:N6	25:RA:2119:A:H62	2.12	0.47
28:RE:3:GLY:HA3	28:RE:81:ILE:HG21	1.97	0.47
28:RE:54:GLN:N	28:RE:54:GLN:CD	2.73	0.47
28:RE:197:ILE:CD1	28:RE:199:ARG:HH12	2.26	0.47
29:RF:33:LEU:HD12	29:RF:33:LEU:HA	1.75	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:RH:53:GLU:CD	31:RH:54:ARG:H	2.21	0.47
36:RQ:86:GLY:O	36:RQ:88:GLY:N	2.46	0.47
39:RT:64:ARG:HD2	39:RT:73:GLU:OE1	2.14	0.47
47:R1:76:ARG:H	47:R1:76:ARG:HD2	1.79	0.47
1:XA:619:U:N3	4:XD:135:LEU:HD23	2.22	0.47
2:XB:93:VAL:HG11	2:XB:97:TRP:CD1	2.50	0.47
3:XC:22:TRP:CD1	3:XC:59:ARG:HD2	2.49	0.47
10:XJ:47:PHE:HB3	14:XN:34:TYR:CE2	2.50	0.47
11:XK:48:ILE:HG13	11:XK:63:LEU:HB2	1.97	0.47
17:XQ:67:LYS:O	17:XQ:68:ARG:HB3	2.15	0.47
20:XT:98:PRO:O	20:XT:100:ILE:N	2.46	0.47
25:YA:213:A:H2'	25:YA:214:G:O4'	2.14	0.47
25:YA:996:A:H4'	40:YU:92:ARG:NE	2.29	0.47
25:YA:1607:C:H4'	25:YA:1608:A:O5'	2.15	0.47
25:YA:2131:G:H1'	25:YA:2158:A:N6	2.28	0.47
25:YA:2378:A:H8	25:YA:2378:A:O5'	1.97	0.47
25:YA:2483:C:H2'	36:YQ:49:ALA:HB1	1.96	0.47
28:YE:52:LEU:HB2	28:YE:75:VAL:CG2	2.40	0.47
28:YE:89:ASP:O	28:YE:90:THR:O	2.33	0.47
28:YE:203:LYS:HD2	28:YE:203:LYS:C	2.39	0.47
30:YG:114:ILE:HD13	30:YG:140:ILE:HG21	1.95	0.47
31:YH:123:PHE:O	31:YH:125:VAL:HG23	2.13	0.47
32:YI:77:LEU:HD22	32:YI:101:LEU:HG	1.96	0.47
36:YQ:59:ARG:CD	36:YQ:59:ARG:N	2.73	0.47
48:Y2:48:HIS:O	48:Y2:49:LYS:C	2.57	0.47
1:QA:9:G:OP2	5:QE:121:LYS:NZ	2.42	0.47
7:QG:18:TYR:HD2	7:QG:59:LEU:HD22	1.79	0.47
25:RA:99:U:H4'	25:RA:101:G:O5'	2.14	0.47
25:RA:221:A:H4'	25:RA:222:A:H5'	1.95	0.47
25:RA:250:G:H2'	25:RA:251:A:C8	2.50	0.47
25:RA:593:G:H1'	54:R8:4:MET:HE1	1.95	0.47
25:RA:981:A:N1	25:RA:2027:G:O2'	2.44	0.47
25:RA:1816:G:H8	27:RD:62:TYR:CZ	2.32	0.47
25:RA:2032:G:H21	28:RE:146:THR:HG23	1.79	0.47
25:RA:2320:A:N3	25:RA:2320:A:H2'	2.29	0.47
25:RA:2344:U:O2'	52:R6:37:ARG:HG2	2.14	0.47
25:RA:2415:G:H4'	35:RP:67:MET:H	1.79	0.47
25:RA:2576:G:O2'	25:RA:2579:C:OP2	2.22	0.47
25:RA:2749:A:H4'	31:RH:62:LYS:HB3	1.96	0.47
27:RD:211:ARG:HD2	27:RD:214:TRP:CZ3	2.50	0.47
28:RE:17:ASP:N	28:RE:17:ASP:OD1	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:RE:65:GLY:HA2	28:RE:70:ALA:HB3	1.95	0.47
30:RG:112:PRO:CG	50:R4:37:SER:HB2	2.45	0.47
36:RQ:66:ILE:H	36:RQ:104:PHE:HA	1.80	0.47
38:RS:56:LEU:O	38:RS:58:LEU:N	2.48	0.47
39:RT:123:GLN:O	39:RT:125:ARG:N	2.48	0.47
42:RW:63:ASP:OD1	42:RW:63:ASP:N	2.48	0.47
50:R4:53:GLU:O	50:R4:57:GLU:HG3	2.14	0.47
54:R8:44:LYS:HD2	54:R8:44:LYS:N	2.30	0.47
1:XA:565:U:H5"	1:XA:566:G:H2'	1.97	0.47
1:XA:1095:U:H5"	1:XA:1109:C:O2	2.14	0.47
1:XA:1203:C:H2'	1:XA:1204:A:O4'	2.14	0.47
9:XI:114:TYR:HD2	10:XJ:60:ARG:HB2	1.79	0.47
15:XO:66:LEU:HD12	15:XO:66:LEU:HA	1.67	0.47
25:YA:28:A:H61	25:YA:512:G:H1'	1.80	0.47
25:YA:181:A:H1'	25:YA:435:C:H5'	1.96	0.47
25:YA:729:G:C5	27:YD:208:LYS:HB2	2.49	0.47
25:YA:1019:U:HO2'	25:YA:1021:A:H2	1.60	0.47
25:YA:1026:U:H1'	25:YA:1027:A:O5'	2.14	0.47
25:YA:1360:A:H2'	25:YA:1361:G:O4'	2.14	0.47
25:YA:1695:G:H1'	27:YD:8:PRO:O	2.14	0.47
25:YA:1930:G:O2'	25:YA:1931:U:P	2.73	0.47
25:YA:2168:G:H8	25:YA:2168:G:OP2	1.97	0.47
25:YA:2512:C:H4'	28:YE:122:PHE:CE2	2.50	0.47
25:YA:2572:A:OP1	28:YE:143:ASN:HB3	2.15	0.47
25:YA:2591:C:H2'	25:YA:2592:G:C8	2.50	0.47
25:YA:2733:A:N1	28:YE:203:LYS:O	2.48	0.47
25:YA:2758:A:C2	25:YA:2759:G:H1'	2.50	0.47
27:YD:32:SER:O	27:YD:33:LEU:CB	2.60	0.47
27:YD:134:ARG:HB2	27:YD:135:PHE:CD1	2.50	0.47
30:YG:28:VAL:O	30:YG:31:VAL:HG13	2.14	0.47
30:YG:34:LEU:HD22	30:YG:35:GLU:N	2.30	0.47
31:YH:67:LEU:O	31:YH:71:LEU:HB2	2.15	0.47
31:YH:104:GLU:HG3	31:YH:114:VAL:HG22	1.96	0.47
33:YN:30:ILE:HG22	33:YN:34:LEU:HD22	1.96	0.47
35:YP:19:VAL:HG12	35:YP:27:HIS:HB2	1.95	0.47
36:YQ:86:GLY:O	36:YQ:88:GLY:N	2.46	0.47
38:YS:19:LYS:O	38:YS:20:ARG:CB	2.55	0.47
1:QA:452:A:H2'	1:QA:453:A:C8	2.50	0.47
1:QA:474:G:H2'	1:QA:475:G:H8	1.80	0.47
1:QA:581:G:OP1	15:QO:61:GLY:HA3	2.14	0.47
1:QA:792:A:H1'	1:QA:793:U:OP2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:1014:A:H4'	19:QS:14:HIS:NE2	2.30	0.47
1:QA:1127:G:H21	1:QA:1147:C:H41	1.63	0.47
1:QA:1287:A:H2	1:QA:1353:G:H1'	1.80	0.47
1:QA:1340:A:HO2'	22:QV:31:G:HO2'	1.61	0.47
2:QB:32:ILE:HD13	2:QB:40:HIS:HB3	1.96	0.47
3:QC:162:GLN:HG2	23:QX:23:A:H2	1.80	0.47
10:QJ:55:LYS:HG3	10:QJ:56:HIS:H	1.78	0.47
19:QS:35:SER:O	19:QS:71:LEU:HD12	2.15	0.47
20:QT:75:ASN:OD1	20:QT:75:ASN:N	2.40	0.47
25:RA:275:G:H3'	25:RA:276:A:H5"	1.97	0.47
25:RA:503:A:H4'	25:RA:504:U:C5'	2.44	0.47
25:RA:1101:U:H2'	25:RA:1102:C:C6	2.49	0.47
25:RA:1278:A:H2'	25:RA:1279:G:C8	2.50	0.47
25:RA:2343:C:O2'	25:RA:2373:G:O2'	2.17	0.47
27:RD:118:VAL:HG22	27:RD:119:ALA:N	2.29	0.47
28:RE:121:ASN:O	28:RE:122:PHE:C	2.57	0.47
35:RP:61:ARG:CD	54:R8:13:ARG:HD2	2.44	0.47
36:RQ:132:VAL:HG11	45:RZ:81:ARG:CZ	2.45	0.47
50:R4:61:ARG:C	50:R4:63:TYR:N	2.73	0.47
1:XA:986:A:H2'	1:XA:987:G:O4'	2.14	0.47
1:XA:1320:C:H42	19:XS:36:ARG:HG3	1.80	0.47
2:XB:18:GLY:H	2:XB:42:ILE:HG22	1.80	0.47
10:XJ:55:LYS:CD	10:XJ:56:HIS:CD2	2.98	0.47
12:XL:120:TYR:O	12:XL:121:GLY:C	2.57	0.47
25:YA:443:A:N7	29:YF:45:ARG:HD2	2.29	0.47
25:YA:2097:C:H2'	25:YA:2098:U:O4'	2.14	0.47
27:YD:72:LYS:CG	27:YD:103:ARG:NH2	2.76	0.47
27:YD:136:ILE:HD12	27:YD:136:ILE:N	2.30	0.47
27:YD:177:LEU:C	27:YD:179:SER:H	2.23	0.47
28:YE:22:PRO:O	28:YE:22:PRO:CG	2.63	0.47
28:YE:54:GLN:N	28:YE:54:GLN:CD	2.73	0.47
28:YE:78:LEU:CD2	28:YE:79:ARG:HD2	2.43	0.47
29:YF:109:GLY:O	29:YF:110:LEU:C	2.58	0.47
29:YF:127:GLU:OE1	29:YF:127:GLU:HA	2.07	0.47
29:YF:162:LEU:HD23	29:YF:165:ARG:HH21	1.79	0.47
30:YG:11:TYR:HA	30:YG:15:VAL:HB	1.95	0.47
32:YI:67:ARG:HH21	32:YI:68:LEU:HB2	1.79	0.47
35:YP:135:LEU:O	35:YP:139:LYS:HB2	2.15	0.47
36:YQ:34:LEU:HD11	36:YQ:129:THR:CB	2.35	0.47
38:YS:46:VAL:HG12	38:YS:47:THR:N	2.28	0.47
49:Y3:23:LEU:HD13	49:Y3:50:VAL:HG11	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:51:A:C6	1:QA:353:A:C2	3.03	0.47
1:QA:262:A:C6	1:QA:263:A:C6	3.03	0.47
1:QA:412:A:H4'	1:QA:413:G:O5'	2.15	0.47
1:QA:1347:G:O2'	1:QA:1348:U:P	2.73	0.47
1:QA:1494:G:OP1	58:QA:1696:PAR:N32	2.48	0.47
3:QC:82:GLU:O	3:QC:86:VAL:HG13	2.14	0.47
11:QK:16:SER:OG	11:QK:106:LYS:NZ	2.48	0.47
25:RA:137(A):G:H2'	25:RA:139:G:N7	2.29	0.47
25:RA:747:U:C1'	51:R5:2:ALA:HB3	2.45	0.47
25:RA:1025:G:C4	25:RA:1135:C:H1'	2.50	0.47
25:RA:1032:A:N1	25:RA:1122:G:O6	2.48	0.47
25:RA:1805:U:O2	27:RD:50:THR:HB	2.15	0.47
25:RA:2291:U:H2'	25:RA:2292:C:C6	2.49	0.47
25:RA:2495:G:H5"	36:RQ:81:VAL:HG12	1.97	0.47
25:RA:2506:U:H1'	56:Z5:101:PPU:HN'3	1.79	0.47
27:RD:12:SER:O	27:RD:16:MET:HB2	2.14	0.47
28:RE:63:LEU:O	28:RE:64:LYS:CB	2.62	0.47
28:RE:101:ARG:C	28:RE:201:THR:OG1	2.58	0.47
31:RH:18:GLU:HA	31:RH:18:GLU:OE2	2.14	0.47
31:RH:41:MET:HG3	31:RH:54:ARG:HA	1.96	0.47
31:RH:126:PRO:CG	31:RH:127:GLU:N	2.65	0.47
40:RU:66:ASN:CG	40:RU:70:ARG:HH21	2.21	0.47
41:RV:24:LYS:HG3	41:RV:92:THR:HG23	1.97	0.47
44:RY:76:CYS:HB2	44:RY:101:LYS:HG3	1.96	0.47
1:XA:1321:C:H5"	1:XA:1322:C:C5'	2.44	0.47
19:XS:41:VAL:HG23	19:XS:67:VAL:HG13	1.97	0.47
25:YA:528:A:C2'	25:YA:529:A:H5'	2.45	0.47
25:YA:760:G:H2'	25:YA:761:A:O4'	2.15	0.47
46:Y0:19:LYS:HA	46:Y0:19:LYS:HD3	1.66	0.47
1:QA:452:A:HO2'	1:QA:453:A:C4'	2.28	0.47
3:QC:131:ARG:HH11	5:QE:50:GLU:HG2	1.80	0.47
25:RA:242:G:H3'	54:R8:6:THR:HG23	1.97	0.47
25:RA:631:A:N3	25:RA:2415:G:O2'	2.41	0.47
25:RA:974(A):C:H4'	25:RA:975:G:O5'	2.15	0.47
25:RA:2439:A:O2'	25:RA:2440:C:OP2	2.26	0.47
25:RA:2681:C:H5'	28:RE:11:MET:SD	2.55	0.47
25:RA:2690:C:OP2	37:RR:14:SER:HB3	2.15	0.47
25:RA:2712:U:OP1	25:RA:2714:G:H4'	2.14	0.47
26:RB:88:C:H2'	26:RB:89:G:O4'	2.15	0.47
36:RQ:34:LEU:HD11	36:RQ:129:THR:CB	2.35	0.47
38:RS:48:LEU:HD23	38:RS:82:ILE:HD11	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:RT:16:ARG:HE	39:RT:19:LEU:HD21	1.80	0.47
40:RU:8:VAL:HG23	40:RU:11:ARG:HH21	1.80	0.47
51:R5:39:MET:C	51:R5:40:LYS:HG3	2.39	0.47
54:R8:56:GLU:C	54:R8:58:ILE:N	2.73	0.47
1:XA:1343:G:H4'	9:XI:122:ALA:HB3	1.97	0.47
5:XE:89:ILE:HG12	5:XE:91:LEU:HD13	1.97	0.47
12:XL:127:GLU:O	12:XL:128:ALA:CB	2.62	0.47
19:XS:41:VAL:HA	19:XS:44:MET:HG3	1.97	0.47
25:YA:2392:A:C8	35:YP:60:MET:HG2	2.50	0.47
25:YA:2404:C:O3'	35:YP:77:ARG:NH2	2.48	0.47
25:YA:2455:G:H2'	25:YA:2456:C:C6	2.50	0.47
25:YA:2600:A:C8	27:YD:237:GLU:OE2	2.67	0.47
27:YD:145:VAL:O	27:YD:153:ALA:HA	2.14	0.47
27:YD:231:HIS:ND1	27:YD:232:PRO:HD2	2.30	0.47
28:YE:15:PHE:CD1	39:YT:81:PRO:HD2	2.50	0.47
28:YE:61:ARG:O	28:YE:63:LEU:CG	2.57	0.47
28:YE:103:ASP:OD2	28:YE:168:MET:HG2	2.15	0.47
28:YE:120:TRP:O	28:YE:121:ASN:HB2	2.15	0.47
31:YH:9:ILE:O	31:YH:10:PRO:O	2.33	0.47
31:YH:18:GLU:HA	31:YH:18:GLU:OE2	2.15	0.47
31:YH:94:TYR:N	31:YH:94:TYR:CD1	2.82	0.47
38:YS:24:LEU:HB2	38:YS:85:VAL:HG12	1.97	0.47
38:YS:28:VAL:HG11	38:YS:98:VAL:HG12	1.97	0.47
38:YS:95:HIS:O	38:YS:96:GLY:C	2.58	0.47
44:YY:73:ARG:HE	44:YY:73:ARG:HB3	1.47	0.47
44:YY:87:LYS:HD3	44:YY:92:ASN:HB3	1.97	0.47
1:QA:222:U:H2'	1:QA:223:U:C6	2.50	0.46
1:QA:224:C:H2'	1:QA:225:C:C6	2.50	0.46
1:QA:1238:A:H62	1:QA:1299:A:N6	2.13	0.46
1:QA:1347:G:H22	1:QA:1374:A:P	2.35	0.46
2:QB:104:ASN:OD1	2:QB:107:THR:OG1	2.30	0.46
10:QJ:81:THR:C	10:QJ:83:GLU:H	2.23	0.46
19:QS:15:LEU:H	19:QS:15:LEU:HD23	1.79	0.46
25:RA:274:G:H1'	25:RA:363:G:C2	2.50	0.46
25:RA:627:A:H4'	25:RA:628:G:H5'	1.97	0.46
25:RA:747:U:O2	25:RA:2014:A:H1'	2.15	0.46
25:RA:2053:G:O6	25:RA:2614:A:H2	1.99	0.46
25:RA:2287:A:N1	25:RA:2346:A:H2	2.13	0.46
25:RA:2420:C:OP1	54:R8:34:TRP:CA	2.63	0.46
25:RA:2593:U:H2'	25:RA:2594:C:H6	1.79	0.46
25:RA:2633:G:H1'	28:RE:62:PRO:HG2	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:RB:89(A):A:C5	26:RB:90:C:H1'	2.50	0.46
27:RD:61:LEU:HA	27:RD:61:LEU:HD12	1.77	0.46
30:RG:105:LYS:HE2	50:R4:26:SER:HB3	1.96	0.46
32:RI:128:LEU:HD13	32:RI:128:LEU:HA	1.77	0.46
34:RO:76:ALA:HB3	39:RT:75:ILE:HB	1.97	0.46
42:RW:23:LEU:O	42:RW:27:LYS:HD2	2.14	0.46
50:R4:3:GLU:HG3	50:R4:4:GLY:H	1.79	0.46
1:XA:545:C:H2'	1:XA:546:G:O4'	2.15	0.46
1:XA:1161:C:O2'	1:XA:1162:C:H5'	2.14	0.46
2:XB:201:ILE:HG21	2:XB:214:ILE:HG21	1.95	0.46
19:XS:33:THR:OG1	19:XS:34:TRP:N	2.49	0.46
24:XY:30:C:N4	24:XY:40:G:H1	2.11	0.46
25:YA:287:C:H2'	25:YA:288:C:C6	2.50	0.46
25:YA:507:A:C5'	25:YA:508:G:H5'	2.44	0.46
25:YA:524:U:H2'	25:YA:525:U:C6	2.50	0.46
25:YA:588:U:C2	29:YF:90:PHE:CE1	3.03	0.46
25:YA:1465:G:H5'	25:YA:1528:A:H1'	1.96	0.46
25:YA:2308:G:C6	25:YA:2311:A:N1	2.83	0.46
25:YA:2345:G:N3	25:YA:2381:C:H2'	2.30	0.46
25:YA:2667:C:N3	31:YH:110:SER:OG	2.44	0.46
25:YA:2729:G:O2'	28:YE:187:ALA:CB	2.63	0.46
28:YE:101:ARG:HD2	28:YE:171:GLU:HA	1.97	0.46
31:YH:86:GLU:O	31:YH:132:ARG:HA	2.16	0.46
31:YH:89:ILE:HD13	31:YH:89:ILE:H	1.80	0.46
36:YQ:87:LYS:O	36:YQ:89:ASN:N	2.43	0.46
38:YS:99:LYS:C	38:YS:101:LEU:N	2.72	0.46
45:YZ:128:VAL:HB	45:YZ:161:VAL:HG13	1.96	0.46
50:Y4:38:LYS:HD3	50:Y4:42:PHE:HE1	1.80	0.46
51:Y5:33:CYS:SG	51:Y5:34:PRO:HD2	2.55	0.46
1:QA:62:U:H3	1:QA:105:G:H1	1.62	0.46
1:QA:300:A:H8	1:QA:300:A:O5'	1.98	0.46
1:QA:565:U:OP2	1:QA:566:G:O2'	2.30	0.46
1:QA:954:G:H21	1:QA:1227:A:H62	1.63	0.46
1:QA:1129:C:O2	1:QA:1132:C:N4	2.39	0.46
2:QB:166:ASP:OD2	2:QB:169:LYS:HB2	2.15	0.46
4:QD:64:LEU:HB2	4:QD:198:VAL:HG11	1.96	0.46
4:QD:187:ARG:NH2	4:QD:190:ASP:HB2	2.31	0.46
14:QN:24:CYS:HB3	14:QN:29:ARG:H	1.79	0.46
17:QQ:74:LEU:HB3	17:QQ:75:ARG:H	1.63	0.46
18:QR:53:ARG:HE	18:QR:59:SER:C	2.22	0.46
19:QS:63:THR:HG23	19:QS:65:ASN:OD1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:QT:89:ARG:NH2	20:QT:105:SER:O	2.37	0.46
25:RA:1535:U:H2'	25:RA:1536:A:C8	2.50	0.46
25:RA:2322:A:H2'	25:RA:2323:G:O4'	2.15	0.46
27:RD:34:VAL:C	27:RD:35:LYS:HG3	2.40	0.46
28:RE:87:GLU:O	28:RE:89:ASP:N	2.49	0.46
28:RE:188:VAL:HG13	28:RE:188:VAL:O	2.15	0.46
29:RF:192:LEU:HD22	29:RF:194:MET:HG2	1.97	0.46
30:RG:3:LEU:HG	50:R4:25:TYR:CE1	2.48	0.46
31:RH:9:ILE:O	31:RH:10:PRO:O	2.33	0.46
31:RH:153:LYS:HG3	31:RH:162:ILE:H	1.78	0.46
39:RT:20:PRO:HD2	39:RT:86:ILE:HG23	1.97	0.46
43:RX:49:VAL:HG13	43:RX:83:VAL:HG13	1.95	0.46
47:R1:89:GLU:HA	47:R1:93:GLU:HB2	1.96	0.46
51:R5:56:LYS:O	51:R5:57:VAL:C	2.57	0.46
1:XA:109:A:H2'	1:XA:326:G:N2	2.30	0.46
3:XC:178:LEU:HD13	3:XC:178:LEU:HA	1.84	0.46
9:XI:83:ARG:O	9:XI:86:VAL:HG12	2.15	0.46
17:XQ:63:ARG:HG2	17:XQ:64:PRO:HD2	1.97	0.46
25:YA:606:U:H4'	25:YA:658:C:H4'	1.97	0.46
25:YA:1174:A:H2'	25:YA:1174:A:N3	2.30	0.46
25:YA:1535:U:C2	25:YA:1537:C:H1'	2.50	0.46
25:YA:2140:C:H2'	25:YA:2141:G:H8	1.80	0.46
25:YA:2564:A:C2	25:YA:2647:U:H4'	2.50	0.46
25:YA:2681:C:H5'	28:YE:11:MET:SD	2.55	0.46
28:YE:188:VAL:O	28:YE:188:VAL:HG13	2.15	0.46
33:YN:114:ARG:O	33:YN:115:ARG:HB3	2.15	0.46
54:Y8:9:GLY:O	54:Y8:13:ARG:HG2	2.16	0.46
1:QA:833:U:H2'	1:QA:834:C:C6	2.50	0.46
2:QB:8:LYS:H	2:QB:8:LYS:HD3	1.79	0.46
5:QE:79:GLU:OE2	8:QH:104:ARG:HA	2.15	0.46
7:QG:78:ARG:HG3	7:QG:79:ARG:N	2.30	0.46
12:QL:43:VAL:HG23	12:QL:93:LEU:HD22	1.96	0.46
16:QP:3:LYS:O	16:QP:21:VAL:HA	2.15	0.46
22:QV:9:G:N2	22:QV:26:G:H1'	2.30	0.46
25:RA:49:A:N7	25:RA:120:U:O4	2.48	0.46
25:RA:49:A:H5"	25:RA:51:G:O4'	2.16	0.46
25:RA:551:G:H5'	25:RA:1220:A:H1'	1.97	0.46
25:RA:576:U:OP1	25:RA:2503:A:OP1	2.32	0.46
25:RA:1051:G:H2'	25:RA:1052:C:O4'	2.15	0.46
25:RA:1203:G:O6	25:RA:1204:A:N6	2.48	0.46
25:RA:1411:C:H42	25:RA:1591:G:H1	1.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2811:G:OP1	28:RE:61:ARG:HG2	2.14	0.46
27:RD:70:TRP:CD2	27:RD:150:LYS:HD2	2.49	0.46
28:RE:33:VAL:HG12	28:RE:90:THR:H	1.80	0.46
30:RG:22:ARG:HH22	30:RG:175:LEU:HD21	1.79	0.46
31:RH:127:GLU:OE2	31:RH:130:ARG:NH2	2.47	0.46
43:RX:55:ASN:HB2	43:RX:80:ILE:HG23	1.97	0.46
50:R4:56:VAL:HA	50:R4:60:GLN:CB	2.28	0.46
1:XA:60:A:P	1:XA:60:A:H8	2.39	0.46
1:XA:407:G:H2'	1:XA:408:A:H8	1.80	0.46
1:XA:562:C:O2'	12:XL:16:GLU:O	2.27	0.46
1:XA:715:A:H2'	1:XA:716:A:C8	2.50	0.46
1:XA:1349:A:H2'	1:XA:1350:A:O4'	2.16	0.46
4:XD:63:LYS:HD2	4:XD:198:VAL:HG22	1.97	0.46
9:XI:18:PHE:HB2	9:XI:62:TYR:HB3	1.97	0.46
19:XS:41:VAL:HB	19:XS:42:PRO:HA	1.96	0.46
25:YA:451:C:H4'	29:YF:52:LYS:HZ2	1.80	0.46
25:YA:2277:G:P	36:YQ:85:LYS:HB2	2.55	0.46
29:YF:46:ARG:CG	29:YF:46:ARG:NH1	2.71	0.46
29:YF:183:VAL:O	29:YF:184:TYR:C	2.57	0.46
31:YH:4:ILE:HG13	31:YH:6:ARG:HD3	1.97	0.46
31:YH:106:THR:HG22	31:YH:112:PRO:HB3	1.97	0.46
36:YQ:109:VAL:HG13	36:YQ:113:GLN:OE1	2.16	0.46
46:Y0:17:GLN:O	46:Y0:19:LYS:HE3	2.14	0.46
54:Y8:29:LYS:HE3	54:Y8:41:ILE:O	2.15	0.46
54:Y8:40:GLU:O	54:Y8:41:ILE:C	2.56	0.46
1:QA:737:A:H2'	1:QA:738:C:C6	2.50	0.46
1:QA:983:A:H1'	1:QA:1049:U:O2	2.15	0.46
1:QA:1402:C:H2'	1:QA:1403:C:O4'	2.15	0.46
6:QF:10:LEU:N	6:QF:59:TYR:O	2.46	0.46
20:QT:53:LEU:HD12	20:QT:100:ILE:HG23	1.98	0.46
25:RA:227:A:C2	25:RA:2407:G:H1'	2.51	0.46
25:RA:928:G:H2'	25:RA:929:G:O4'	2.16	0.46
25:RA:2477:C:H3'	55:R9:1:MET:HE3	1.96	0.46
27:RD:35:LYS:HZ1	27:RD:65:ILE:HA	1.80	0.46
28:RE:94:GLU:C	28:RE:96:PHE:N	2.73	0.46
31:RH:4:ILE:H	31:RH:4:ILE:CD1	2.25	0.46
44:RY:89:PHE:O	44:RY:90:LEU:HD13	2.15	0.46
51:R5:43:HIS:N	51:R5:43:HIS:ND1	2.63	0.46
54:R8:29:LYS:HE3	54:R8:41:ILE:O	2.15	0.46
1:XA:335:C:H2'	1:XA:336:C:C6	2.51	0.46
1:XA:345:C:H4'	1:XA:346:G:O5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:963:G:H21	10:XJ:54:PHE:HE1	1.61	0.46
1:XA:1062:U:H2'	1:XA:1063:C:C6	2.50	0.46
1:XA:1250:A:H2'	1:XA:1251:A:C8	2.50	0.46
1:XA:1319:A:O2'	1:XA:1323:G:N7	2.39	0.46
3:XC:14:ILE:HG12	3:XC:15:THR:N	2.30	0.46
7:XG:115:ARG:HB2	7:XG:118:VAL:HG22	1.97	0.46
10:XJ:55:LYS:CG	10:XJ:56:HIS:H	2.27	0.46
16:XP:1:MET:O	16:XP:3:LYS:HG3	2.15	0.46
25:YA:964:C:O2'	25:YA:2273:A:N3	2.43	0.46
25:YA:2783:G:H2'	25:YA:2784:C:C6	2.51	0.46
27:YD:61:LEU:HB3	27:YD:63:ARG:NH1	2.31	0.46
27:YD:110:GLY:O	27:YD:111:LEU:C	2.58	0.46
28:YE:3:GLY:HA3	28:YE:81:ILE:HG21	1.97	0.46
32:YI:97:ILE:HD12	32:YI:140:LEU:HD11	1.97	0.46
36:YQ:63:LYS:HE2	36:YQ:65:PHE:CZ	2.50	0.46
50:Y4:2:LYS:HA	50:Y4:2:LYS:HD2	1.67	0.46
1:QA:689:C:OP1	11:QK:27:ASN:ND2	2.49	0.46
1:QA:715:A:H2'	1:QA:716:A:C8	2.50	0.46
1:QA:1225:A:H2'	1:QA:1225:A:N3	2.31	0.46
12:QL:115:LYS:O	12:QL:117:ARG:N	2.47	0.46
25:RA:945:A:C4	25:RA:2448:A:C2	3.03	0.46
25:RA:1918:A:O2'	25:RA:1920:C:N4	2.48	0.46
25:RA:2015:A:H1'	51:R5:2:ALA:CA	2.25	0.46
28:RE:54:GLN:CA	28:RE:54:GLN:HE21	2.27	0.46
28:RE:89:ASP:O	28:RE:90:THR:O	2.33	0.46
31:RH:86:GLU:O	31:RH:132:ARG:HA	2.15	0.46
38:RS:61:ASN:O	38:RS:65:VAL:HG23	2.14	0.46
38:RS:78:LEU:HD11	38:RS:107:GLU:O	2.15	0.46
42:RW:86:LEU:O	42:RW:94:ASP:N	2.44	0.46
43:RX:40:LYS:C	43:RX:42:ALA:H	2.23	0.46
50:R4:64:GLY:C	50:R4:66:SER:N	2.73	0.46
54:R8:52:LYS:O	54:R8:52:LYS:CG	2.64	0.46
1:XA:156:G:H1	1:XA:165:C:N4	2.12	0.46
1:XA:323:U:H2'	1:XA:324:G:O4'	2.16	0.46
2:XB:163:PHE:CD1	2:XB:185:ILE:HG13	2.50	0.46
7:XG:74:GLU:HG2	7:XG:91:VAL:HG22	1.98	0.46
8:XH:1:MET:HE2	8:XH:1:MET:H3	1.81	0.46
10:XJ:62:HIS:H	10:XJ:62:HIS:CD2	2.33	0.46
25:YA:573:G:N1	25:YA:2031:A:OP2	2.40	0.46
25:YA:727:A:C2	27:YD:9:TYR:CD2	3.04	0.46
25:YA:877:U:H3	25:YA:899:A:H2	1.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1655:A:O3'	28:YE:115:GLY:HA3	2.15	0.46
25:YA:2820:A:O2'	25:YA:2821:A:OP1	2.32	0.46
27:YD:35:LYS:HE3	27:YD:65:ILE:N	2.31	0.46
27:YD:102:LYS:O	27:YD:103:ARG:HG3	2.15	0.46
27:YD:206:LEU:HA	27:YD:206:LEU:HD23	1.49	0.46
29:YF:184:TYR:CD2	29:YF:188:ARG:HD2	2.50	0.46
32:YI:125:GLU:OE1	32:YI:141:LYS:HB3	2.15	0.46
36:YQ:66:ILE:H	36:YQ:104:PHE:HA	1.79	0.46
38:YS:61:ASN:O	38:YS:65:VAL:HG23	2.15	0.46
40:YU:83:LEU:HG	40:YU:88:ILE:HG13	1.97	0.46
47:Y1:96:LYS:H	47:Y1:97:LEU:HD12	1.81	0.46
48:Y2:17:SER:HB2	48:Y2:18:PRO:HG3	1.98	0.46
1:QA:1376:U:P	7:QG:94:ARG:HH12	2.39	0.46
4:QD:7:PRO:HB2	4:QD:10:ARG:HD2	1.98	0.46
4:QD:30:LYS:C	4:QD:32:ALA:HA	2.41	0.46
25:RA:1312:U:H4'	25:RA:1313:U:O5'	2.16	0.46
25:RA:1651:G:H2'	25:RA:1652:A:O4'	2.16	0.46
25:RA:2146:C:H4'	25:RA:2147:G:C8	2.50	0.46
25:RA:2405:G:O2'	25:RA:2406:U:OP2	2.31	0.46
28:RE:50:GLY:CA	28:RE:74:PRO:HG3	2.45	0.46
31:RH:59:ARG:CG	31:RH:59:ARG:NH1	2.79	0.46
32:RI:88:ILE:HG12	32:RI:122:GLU:H	1.81	0.46
38:RS:24:LEU:HB2	38:RS:85:VAL:HG12	1.96	0.46
52:R6:13:CYS:HB2	52:R6:22:ALA:HB3	1.98	0.46
11:XK:88:GLY:C	11:XK:90:GLY:H	2.23	0.46
15:XO:56:LEU:O	15:XO:60:VAL:HG23	2.16	0.46
18:XR:32:ARG:HA	18:XR:69:THR:HG21	1.97	0.46
25:YA:363(B):G:H2'	25:YA:363(C):G:C8	2.51	0.46
25:YA:1476:C:H2'	25:YA:1477:A:C8	2.51	0.46
25:YA:2361:A:O5'	54:Y8:27:THR:OG1	2.25	0.46
25:YA:2653:U:O2'	31:YH:110:SER:HB2	2.16	0.46
26:YB:91:C:H5"	45:YZ:79:ARG:NH1	2.31	0.46
27:YD:35:LYS:HD3	27:YD:63:ARG:HB3	1.96	0.46
27:YD:48:ARG:HG3	27:YD:48:ARG:NH1	2.31	0.46
27:YD:211:ARG:HG2	27:YD:211:ARG:HH11	1.80	0.46
28:YE:54:GLN:CA	28:YE:54:GLN:HE21	2.27	0.46
31:YH:86:GLU:O	31:YH:87:LEU:CB	2.64	0.46
31:YH:151:ILE:O	31:YH:152:ARG:O	2.34	0.46
35:YP:57:THR:C	35:YP:59:LEU:H	2.24	0.46
38:YS:110:LEU:HA	38:YS:112:PHE:CZ	2.50	0.46
48:Y2:15:LYS:H	48:Y2:67:LYS:HZ3	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:Y3:8:LEU:HD22	49:Y3:31:LEU:HD22	1.96	0.46
51:Y5:48:GLU:HA	51:Y5:59:GLU:HG2	1.97	0.46
54:Y8:44:LYS:HD2	54:Y8:44:LYS:N	2.30	0.46
54:Y8:48:PHE:N	54:Y8:48:PHE:HD1	2.14	0.46
1:QA:1222:G:C2	1:QA:1223:C:C2	3.04	0.46
2:QB:165:VAL:HG23	2:QB:166:ASP:H	1.81	0.46
4:QD:9:CYS:SG	4:QD:32:ALA:N	2.89	0.46
4:QD:78:LEU:HD22	4:QD:96:LEU:HB3	1.98	0.46
5:QE:41:VAL:HG13	5:QE:113:ALA:HA	1.98	0.46
13:QM:84:ILE:HD12	13:QM:84:ILE:HA	1.74	0.46
25:RA:307:G:N2	25:RA:309:G:H3'	2.30	0.46
25:RA:1303:G:H1'	25:RA:1641:A:N1	2.31	0.46
25:RA:2154:G:H2'	25:RA:2155:G:H8	1.81	0.46
25:RA:2469:A:H5"	25:RA:2470:G:C8	2.51	0.46
25:RA:2788:C:O2'	25:RA:2809:A:N3	2.48	0.46
27:RD:35:LYS:HE3	27:RD:63:ARG:C	2.41	0.46
28:RE:103:ASP:OD2	28:RE:168:MET:HG2	2.15	0.46
30:RG:65:GLY:C	50:R4:7:PRO:HG2	2.40	0.46
36:RQ:66:ILE:O	36:RQ:104:PHE:N	2.49	0.46
36:RQ:109:VAL:HG13	36:RQ:113:GLN:OE1	2.16	0.46
38:RS:83:LYS:O	38:RS:109:GLY:HA3	2.15	0.46
1:XA:92:G:H3'	1:XA:93:U:C6	2.51	0.46
1:XA:411:A:N9	1:XA:413:G:H1'	2.31	0.46
1:XA:1314:C:N4	19:XS:2:PRO:O	2.48	0.46
2:XB:70:PHE:O	2:XB:93:VAL:N	2.33	0.46
9:XI:18:PHE:HD2	9:XI:62:TYR:HD2	1.61	0.46
20:XT:45:GLN:HB2	20:XT:91:LEU:HD13	1.96	0.46
25:YA:636:G:OP1	35:YP:132:LYS:HB2	2.15	0.46
25:YA:1203:G:H3'	25:YA:1204:A:H5"	1.97	0.46
25:YA:1264:G:H3'	25:YA:1265:A:H5"	1.97	0.46
25:YA:2667:C:H1'	31:YH:109:PHE:CD2	2.42	0.46
25:YA:2774:C:H2'	25:YA:2775:A:O4'	2.14	0.46
26:YB:24:G:H1'	26:YB:27:C:N4	2.31	0.46
27:YD:79:VAL:HG21	27:YD:111:LEU:HD21	1.98	0.46
27:YD:148:GLU:HB2	27:YD:151:LYS:HD2	1.98	0.46
28:YE:63:LEU:O	28:YE:64:LYS:CB	2.62	0.46
28:YE:188:VAL:HA	28:YE:189:PRO:HD2	1.79	0.46
28:YE:195:LEU:HD12	28:YE:196:VAL:N	2.29	0.46
30:YG:34:LEU:HD12	30:YG:100:TRP:CH2	2.50	0.46
1:QA:1348:U:H3	1:QA:1374:A:H2	1.64	0.46
10:QJ:84:GLN:H	10:QJ:84:GLN:HG3	1.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:QT:29:LYS:O	20:QT:33:ILE:HG12	2.16	0.46
25:RA:483:A:H5"	44:RY:49:VAL:HG13	1.98	0.46
25:RA:586:A:N1	25:RA:809:G:O2'	2.33	0.46
25:RA:631:A:OP1	35:RP:64:LYS:HE2	2.16	0.46
25:RA:676:A:N1	25:RA:802:A:N1	2.64	0.46
25:RA:752:A:H3'	53:R7:1:MET:SD	2.55	0.46
25:RA:845:G:H8	25:RA:845:G:OP2	1.99	0.46
28:RE:61:ARG:O	28:RE:63:LEU:CG	2.57	0.46
28:RE:172:VAL:HG13	28:RE:182:LEU:HD11	1.98	0.46
30:RG:98:ARG:O	30:RG:101:ILE:HG13	2.16	0.46
31:RH:37:VAL:HG11	31:RH:68:THR:HG23	1.98	0.46
32:RI:29:TYR:O	32:RI:33:ARG:HB2	2.16	0.46
32:RI:57:ARG:HA	32:RI:60:GLU:HB3	1.96	0.46
36:RQ:59:ARG:C	36:RQ:60:ARG:CG	2.74	0.46
36:RQ:63:LYS:HD2	45:RZ:175:VAL:HG21	1.98	0.46
44:RY:17:SER:OG	44:RY:71:LYS:HD2	2.16	0.46
45:RZ:74:VAL:HG13	45:RZ:86:VAL:HG22	1.98	0.46
50:R4:4:GLY:O	50:R4:5:ILE:C	2.59	0.46
52:R6:33:LYS:HG3	52:R6:34:LEU:HD13	1.98	0.46
1:XA:222:U:H2'	1:XA:223:U:C6	2.51	0.46
1:XA:620:C:H2'	1:XA:621:A:O4'	2.16	0.46
1:XA:762:C:H2'	1:XA:763:G:C8	2.51	0.46
7:XG:50:ILE:HG21	7:XG:61:VAL:HG21	1.98	0.46
9:XI:46:ALA:HA	9:XI:78:LYS:HB2	1.98	0.46
25:YA:271(B):G:N7	25:YA:421:U:H2'	2.31	0.46
25:YA:466:A:N3	25:YA:683:C:H1'	2.31	0.46
25:YA:1062:G:H8	25:YA:1062:G:O5'	1.98	0.46
25:YA:1408:C:C2	25:YA:1595:G:N2	2.84	0.46
25:YA:2025:C:H2'	25:YA:2026:C:C6	2.51	0.46
25:YA:2150:U:H2'	25:YA:2151:G:C8	2.51	0.46
25:YA:2168:G:N2	25:YA:2170:A:H62	2.14	0.46
25:YA:2310:A:N6	30:YG:79:ASN:HB2	2.30	0.46
27:YD:2:ALA:HB1	27:YD:20:ASP:CB	2.46	0.46
27:YD:36:PRO:HB3	27:YD:62:TYR:O	2.16	0.46
28:YE:87:GLU:O	28:YE:89:ASP:N	2.48	0.46
28:YE:111:ARG:NE	28:YE:160:TYR:CE1	2.76	0.46
28:YE:137:HIS:CB	28:YE:138:PRO:HD2	2.42	0.46
31:YH:88:LEU:HD22	31:YH:163:TYR:O	2.16	0.46
31:YH:128:PRO:HD2	31:YH:129:THR:N	2.25	0.46
33:YN:9:VAL:HG23	33:YN:10:GLU:N	2.31	0.46
36:YQ:87:LYS:O	36:YQ:87:LYS:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:YS:108:GLY:O	38:YS:110:LEU:N	2.48	0.46
39:YT:95:ARG:HD2	39:YT:95:ARG:HA	1.75	0.46
40:YU:75:ASN:HB3	40:YU:78:THR:H	1.81	0.46
47:Y1:80:LEU:HB2	47:Y1:81:LYS:H	1.61	0.46
54:Y8:56:GLU:C	54:Y8:58:ILE:N	2.73	0.46
1:QA:17:U:H2'	1:QA:18:C:C6	2.50	0.46
1:QA:335:C:H2'	1:QA:336:C:C6	2.51	0.46
1:QA:1446:A:HO2'	1:QA:1447:G:P	2.39	0.46
10:QJ:31:GLY:HA3	10:QJ:78:ASN:CG	2.40	0.46
10:QJ:55:LYS:CD	10:QJ:56:HIS:CD2	2.98	0.46
12:QL:27:LEU:HD13	12:QL:28:LYS:N	2.30	0.46
19:QS:32:LYS:HA	19:QS:50:ALA:HB3	1.98	0.46
25:RA:321:G:H5"	29:RF:136:THR:HG23	1.98	0.46
25:RA:1026:U:H1'	25:RA:1027:A:H5"	1.98	0.46
25:RA:1497:U:H5"	25:RA:1498:C:H5	1.80	0.46
27:RD:68:LYS:HD2	27:RD:70:TRP:CZ2	2.51	0.46
31:RH:4:ILE:HG13	31:RH:6:ARG:HD3	1.97	0.46
31:RH:13:LYS:HE2	31:RH:13:LYS:CA	2.40	0.46
31:RH:66:GLY:O	31:RH:67:LEU:C	2.58	0.46
34:RO:31:LYS:HB3	34:RO:32:TYR:CD2	2.51	0.46
36:RQ:26:TYR:O	36:RQ:27:VAL:O	2.34	0.46
36:RQ:30:GLY:CA	36:RQ:107:ALA:HB2	2.39	0.46
36:RQ:63:LYS:HE2	36:RQ:65:PHE:CZ	2.50	0.46
39:RT:26:ASP:HB2	39:RT:90:GLN:O	2.16	0.46
41:RV:51:VAL:HG12	41:RV:53:GLU:H	1.80	0.46
50:R4:26:SER:C	50:R4:27:THR:O	2.58	0.46
55:R9:8:LYS:O	55:R9:34:GLN:NE2	2.49	0.46
1:XA:24:U:H2'	1:XA:25:C:C6	2.51	0.46
1:XA:130:A:N3	1:XA:263:A:O2'	2.47	0.46
1:XA:745:C:OP1	1:XA:851:G:O2'	2.33	0.46
2:XB:114:ARG:O	2:XB:118:LEU:HG	2.16	0.46
3:XC:78:GLY:HA3	3:XC:83:ARG:HB3	1.98	0.46
6:XF:48:LEU:HG	6:XF:57:GLN:HA	1.98	0.46
12:XL:27:LEU:HD13	12:XL:28:LYS:N	2.30	0.46
20:XT:26:ASN:O	20:XT:30:LYS:HB2	2.16	0.46
25:YA:27:G:H1'	25:YA:513:A:N6	2.31	0.46
25:YA:478:A:C6	25:YA:480:A:C6	3.04	0.46
25:YA:954:G:OP1	36:YQ:15:GLY:N	2.43	0.46
25:YA:1657:C:H2'	25:YA:1658:C:C6	2.51	0.46
27:YD:14:ARG:HG3	27:YD:15:PHE:N	2.31	0.46
27:YD:105:ILE:HG23	27:YD:106:ILE:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:183:ARG:CG	27:YD:183:ARG:NH1	2.69	0.46
28:YE:36:ARG:O	28:YE:37:ARG:C	2.59	0.46
36:YQ:11:LYS:HE2	36:YQ:87:LYS:HA	1.98	0.46
38:YS:13:ARG:O	38:YS:14:VAL:HB	2.15	0.46
38:YS:102:ALA:C	38:YS:104:GLY:N	2.73	0.46
42:YW:110:LYS:HG3	42:YW:111:HIS:ND1	2.31	0.46
1:QA:1347:G:HO2'	1:QA:1348:U:P	2.38	0.46
1:QA:1443:G:C6	39:RT:118:ARG:HB2	2.51	0.46
5:QE:75:THR:CB	5:QE:117:ASP:O	2.63	0.46
25:RA:392:C:H5"	25:RA:409:C:H5"	1.97	0.46
25:RA:699:A:H2'	25:RA:700:G:O4'	2.16	0.46
25:RA:861:A:N3	26:RB:79:C:O2'	2.49	0.46
25:RA:1885:A:H5'	25:RA:1886:C:OP2	2.16	0.46
25:RA:2321:G:N3	25:RA:2321:G:C2'	2.77	0.46
25:RA:2335:A:O2'	25:RA:2336:A:H2'	2.16	0.46
28:RE:47:VAL:O	28:RE:47:VAL:HG23	2.16	0.46
29:RF:11:VAL:HG12	29:RF:12:LEU:H	1.80	0.46
29:RF:20:LEU:HD23	29:RF:125:LEU:HD12	1.98	0.46
31:RH:88:LEU:HD22	31:RH:163:TYR:O	2.16	0.46
36:RQ:27:VAL:HG22	36:RQ:105:GLU:CD	2.41	0.46
36:RQ:87:LYS:O	36:RQ:87:LYS:HG2	2.15	0.46
37:RR:37:THR:OG1	37:RR:40:LYS:HG3	2.16	0.46
40:RU:69:CYS:HB3	40:RU:106:PHE:HZ	1.81	0.46
47:R1:73:LEU:HB3	47:R1:90:ILE:HG23	1.97	0.46
51:R5:41:PRO:HA	51:R5:42:PRO:HD3	1.82	0.46
52:R6:44:ARG:O	52:R6:45:LYS:HB2	2.16	0.46
1:XA:926:G:C6	1:XA:1505:G:C6	3.04	0.46
2:XB:217:ARG:HE	2:XB:217:ARG:HB2	1.54	0.46
20:XT:33:ILE:HD13	20:XT:62:LEU:HB3	1.98	0.46
25:YA:1512:G:H2'	25:YA:1513:C:C6	2.50	0.46
25:YA:1655:A:H3'	25:YA:1656:C:H6	1.81	0.46
25:YA:2208:U:C1'	27:YD:151:LYS:HE2	2.45	0.46
27:YD:118:VAL:O	27:YD:129:ASN:HA	2.16	0.46
29:YF:31:HIS:O	29:YF:34:TRP:HB3	2.15	0.46
31:YH:51:ARG:HG3	31:YH:51:ARG:NH1	2.30	0.46
36:YQ:27:VAL:HG22	36:YQ:105:GLU:CD	2.41	0.46
37:YR:51:LEU:HD12	37:YR:70:LEU:HG	1.97	0.46
37:YR:78:LYS:O	37:YR:83:ILE:HG12	2.16	0.46
44:YY:84:ARG:HB3	44:YY:95:LYS:HD3	1.97	0.46
1:QA:1298:C:H4'	1:QA:1299:A:N9	2.31	0.45
2:QB:120:ALA:C	2:QB:122:PHE:H	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:QC:19:GLU:HA	3:QC:54:ARG:HH12	1.81	0.45
6:QF:99:ALA:HB1	18:QR:23:LYS:NZ	2.31	0.45
10:QJ:22:LYS:HB3	10:QJ:22:LYS:HE3	1.68	0.45
25:RA:678:C:H2'	25:RA:679:C:C6	2.51	0.45
25:RA:1153:C:H2'	25:RA:1154:G:O4'	2.16	0.45
25:RA:1301:A:O2'	25:RA:1302:A:H3'	2.16	0.45
25:RA:2469:A:H2	25:RA:2481:G:H21	1.64	0.45
25:RA:2645:G:N2	25:RA:2767:C:OP2	2.49	0.45
29:RF:182:ASN:O	29:RF:186:ILE:HG12	2.16	0.45
32:RI:113:ARG:HG3	32:RI:131:LYS:NZ	2.30	0.45
36:RQ:81:VAL:HG23	36:RQ:82:ARG:N	2.31	0.45
36:RQ:85:LYS:HD3	36:RQ:86:GLY:H	1.80	0.45
37:RR:33:ARG:HG2	37:RR:34:ILE:N	2.30	0.45
37:RR:70:LEU:C	37:RR:72:ASP:H	2.21	0.45
50:R4:15:ILE:HG22	50:R4:20:ASN:CA	2.45	0.45
50:R4:50:VAL:O	50:R4:50:VAL:HG13	2.15	0.45
1:XA:154:C:H42	1:XA:167:G:H1	1.64	0.45
1:XA:626:U:C2	1:XA:627:G:C8	3.05	0.45
10:XJ:55:LYS:CG	10:XJ:56:HIS:CD2	2.97	0.45
16:XP:8:ARG:C	16:XP:9:PHE:HD1	2.24	0.45
18:XR:73:ALA:HB3	18:XR:79:LEU:HD12	1.98	0.45
20:XT:29:LYS:CD	20:XT:66:ALA:HB2	2.45	0.45
25:YA:508:G:HO2'	25:YA:509:C:P	2.39	0.45
25:YA:2016:U:H1'	51:Y5:6:VAL:HG13	1.97	0.45
27:YD:11:PRO:O	27:YD:12:SER:CB	2.65	0.45
27:YD:18:VAL:CG1	27:YD:19:ALA:N	2.78	0.45
28:YE:62:PRO:O	28:YE:63:LEU:C	2.59	0.45
28:YE:94:GLU:C	28:YE:96:PHE:N	2.73	0.45
28:YE:101:ARG:C	28:YE:201:THR:OG1	2.58	0.45
29:YF:65:TRP:CH2	29:YF:72:ARG:HB3	2.50	0.45
31:YH:109:PHE:CE1	31:YH:152:ARG:NH1	2.85	0.45
36:YQ:85:LYS:HD3	36:YQ:86:GLY:H	1.80	0.45
41:YV:19:LYS:HA	41:YV:94:LEU:O	2.15	0.45
47:Y1:58:ILE:HD12	47:Y1:58:ILE:N	2.30	0.45
47:Y1:83:GLU:C	47:Y1:85:LEU:H	2.24	0.45
1:QA:491:G:H2'	1:QA:492:G:O4'	2.16	0.45
1:QA:757:U:H2'	1:QA:758:G:O4'	2.16	0.45
1:QA:1378:C:N4	1:QA:1379:G:N3	2.64	0.45
3:QC:23:TYR:CD1	10:QJ:10:GLY:HA2	2.51	0.45
4:QD:46:LYS:HD3	4:QD:46:LYS:HA	1.69	0.45
8:QH:20:TYR:HA	8:QH:65:TYR:CZ	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:QI:45:ALA:O	9:QI:48:GLU:HG2	2.15	0.45
25:RA:297:C:H5"	44:RY:85:VAL:HG21	1.96	0.45
25:RA:957:A:H5'	36:RQ:76:LYS:CD	2.47	0.45
25:RA:1939:U:OP1	25:RA:2604:U:O2'	2.33	0.45
25:RA:2131:G:N2	25:RA:2158:A:N7	2.63	0.45
25:RA:2516:G:C6	25:RA:2517:C:N4	2.84	0.45
25:RA:2698:U:H2'	25:RA:2699:C:C6	2.51	0.45
26:RB:5:C:O2'	26:RB:27:C:O2	2.33	0.45
28:RE:1:MET:HE2	28:RE:1:MET:N	2.31	0.45
28:RE:15:PHE:CD1	28:RE:20:ALA:HB2	2.50	0.45
31:RH:89:ILE:H	31:RH:89:ILE:HD13	1.81	0.45
32:RI:52:ARG:CB	32:RI:56:LYS:HB3	2.45	0.45
33:RN:57:ALA:C	33:RN:60:ILE:HD11	2.40	0.45
2:XB:55:PHE:HD1	2:XB:58:ILE:HG13	1.81	0.45
25:YA:1263:U:H1'	51:Y5:10:LYS:HG3	1.98	0.45
25:YA:1683:C:H2'	25:YA:1684:C:C6	2.51	0.45
25:YA:2619:C:H1'	28:YE:156:MET:HE1	1.97	0.45
28:YE:1:MET:HA	28:YE:200:GLU:OE2	2.16	0.45
28:YE:1:MET:HE2	28:YE:1:MET:N	2.31	0.45
28:YE:77:ILE:O	28:YE:78:LEU:O	2.35	0.45
29:YF:7:TYR:N	29:YF:7:TYR:CD1	2.84	0.45
29:YF:167:ALA:HB1	29:YF:173:VAL:HG11	1.98	0.45
35:YP:115:LEU:HB3	35:YP:131:SER:HB2	1.99	0.45
37:YR:101:ALA:CB	51:Y5:47:PRO:HD3	2.46	0.45
44:YY:87:LYS:HA	44:YY:92:ASN:HB3	1.98	0.45
52:Y6:33:LYS:HE2	52:Y6:33:LYS:HB2	1.76	0.45
53:Y7:22:MET:HE2	53:Y7:31:LEU:CD1	2.46	0.45
1:QA:662:G:H2'	1:QA:663:A:C8	2.51	0.45
1:QA:695:A:H2'	1:QA:696:A:C8	2.51	0.45
1:QA:1014:A:H2'	1:QA:1015:A:C8	2.52	0.45
7:QG:20:ASP:OD1	7:QG:21:VAL:N	2.49	0.45
7:QG:45:ASP:O	7:QG:49:ILE:HG12	2.16	0.45
10:QJ:55:LYS:HG3	10:QJ:56:HIS:CG	2.51	0.45
10:QJ:58:ASP:O	10:QJ:59:SER:HB2	2.17	0.45
12:QL:61:THR:O	12:QL:63:GLY:N	2.45	0.45
12:QL:113:ARG:NH2	12:QL:120:TYR:CE2	2.85	0.45
25:RA:307:G:O5'	25:RA:307:G:H8	1.99	0.45
25:RA:1796:U:H2'	25:RA:1797:C:H6	1.79	0.45
25:RA:2021:C:OP1	51:R5:12:SER:OG	2.25	0.45
25:RA:2287:A:N1	25:RA:2346:A:C2	2.84	0.45
25:RA:2290:G:H1	25:RA:2342:C:H42	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:RB:49:C:H2'	26:RB:50:G:C8	2.51	0.45
28:RE:1:MET:HA	28:RE:200:GLU:OE2	2.16	0.45
28:RE:62:PRO:O	28:RE:63:LEU:C	2.59	0.45
30:RG:106:LEU:HA	30:RG:110:ALA:HB3	1.98	0.45
31:RH:137:ASP:HB2	31:RH:140:LYS:HE3	1.98	0.45
31:RH:151:ILE:O	31:RH:152:ARG:O	2.34	0.45
36:RQ:52:VAL:O	36:RQ:53:ALA:C	2.59	0.45
36:RQ:104:PHE:O	36:RQ:105:GLU:CB	2.65	0.45
1:XA:190:G:HO2'	1:XA:191(A):G:P	2.40	0.45
1:XA:474:G:H2'	1:XA:475:G:C8	2.51	0.45
1:XA:1252:A:H2'	1:XA:1253:G:O4'	2.16	0.45
1:XA:1392:G:H2'	1:XA:1393:U:C6	2.51	0.45
8:XH:91:ARG:HG3	12:XL:7:ILE:HG21	1.98	0.45
10:XJ:50:ILE:HD11	10:XJ:57:LYS:CD	2.46	0.45
12:XL:115:LYS:O	12:XL:117:ARG:N	2.47	0.45
25:YA:34:C:H41	25:YA:447:A:H61	1.65	0.45
25:YA:975:G:H1'	25:YA:990:A:C2	2.51	0.45
25:YA:1054:A:N1	25:YA:1105:U:O4	2.49	0.45
25:YA:1316:U:H2'	25:YA:1317:A:H8	1.81	0.45
25:YA:1728:G:H8	25:YA:1732:A:H62	1.64	0.45
25:YA:2334:G:C6	46:Y0:75:LEU:HD21	2.52	0.45
25:YA:2347:C:H2'	25:YA:2348:U:C6	2.51	0.45
25:YA:2867:G:OP2	39:YT:119:LYS:NZ	2.28	0.45
27:YD:65:ILE:HD11	27:YD:67:PHE:CE2	2.51	0.45
28:YE:51:PHE:HD2	28:YE:52:LEU:H	1.59	0.45
29:YF:117:ARG:NH2	29:YF:189:THR:O	2.49	0.45
30:YG:146:TYR:O	30:YG:149:VAL:HG22	2.16	0.45
31:YH:16:SER:OG	31:YH:17:VAL:N	2.50	0.45
37:YR:24:GLN:HE21	37:YR:44:LEU:HG	1.81	0.45
37:YR:109:ALA:HA	37:YR:110:PRO:HD2	1.77	0.45
38:YS:109:GLY:O	38:YS:110:LEU:HB2	2.16	0.45
40:YU:58:ARG:HA	40:YU:61:TRP:CE3	2.51	0.45
40:YU:66:ASN:O	40:YU:70:ARG:HB2	2.17	0.45
40:YU:68:ALA:O	40:YU:71:GLN:HB2	2.16	0.45
44:YY:51:VAL:HG13	44:YY:52:SER:N	2.28	0.45
44:YY:56:PRO:O	44:YY:58:GLY:N	2.49	0.45
45:YZ:149:SER:HB2	45:YZ:172:ALA:O	2.16	0.45
48:Y2:17:SER:HB2	48:Y2:18:PRO:CG	2.46	0.45
1:QA:49:U:O4	1:QA:365:U:O4	2.34	0.45
1:QA:106:C:H2'	1:QA:107:G:H8	1.80	0.45
1:QA:280:C:H3'	1:QA:281:G:H5'	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:390:C:O3'	16:QP:28:ARG:NH2	2.44	0.45
1:QA:1120:G:H2'	1:QA:1121:U:C6	2.51	0.45
1:QA:1367:C:H4'	10:QJ:48:THR:HG21	1.98	0.45
4:QD:98:GLU:OE2	4:QD:107:ARG:NE	2.49	0.45
12:QL:64:TYR:O	12:QL:65:GLU:HB2	2.16	0.45
25:RA:2335:A:O2'	25:RA:2336:A:H8	1.99	0.45
25:RA:2405:G:O2'	25:RA:2406:U:P	2.74	0.45
25:RA:2655:G:O2'	25:RA:2656:U:H5	1.98	0.45
27:RD:30:GLU:HG3	27:RD:63:ARG:CZ	2.47	0.45
27:RD:121:PRO:HB3	27:RD:135:PHE:CE2	2.52	0.45
28:RE:21:VAL:HG23	28:RE:22:PRO:CD	2.46	0.45
29:RF:65:TRP:O	29:RF:67:GLN:N	2.42	0.45
31:RH:106:THR:HG22	31:RH:112:PRO:HB3	1.97	0.45
31:RH:109:PHE:CE1	31:RH:152:ARG:NH1	2.84	0.45
36:RQ:5:ARG:O	36:RQ:6:ARG:O	2.35	0.45
38:RS:16:ASN:HA	38:RS:19:LYS:HD3	1.98	0.45
51:R5:16:ARG:O	51:R5:20:ARG:HG3	2.16	0.45
52:R6:26:ASN:ND2	52:R6:35:GLU:OE2	2.49	0.45
1:XA:19:C:H2'	1:XA:20:U:H6	1.81	0.45
2:XB:140:HIS:HA	2:XB:143:GLU:OE1	2.17	0.45
7:XG:15:ASP:OD1	7:XG:44:TYR:OH	2.35	0.45
8:XH:86:ILE:HG13	8:XH:133:LEU:HD22	1.98	0.45
12:XL:10:LEU:HB3	17:XQ:32:TYR:CE2	2.52	0.45
12:XL:64:TYR:O	12:XL:65:GLU:HB2	2.16	0.45
12:XL:113:ARG:NH2	12:XL:120:TYR:CE2	2.84	0.45
13:XM:58:GLU:O	13:XM:62:ASN:ND2	2.33	0.45
17:XQ:43:LEU:HD12	17:XQ:68:ARG:HG2	1.97	0.45
20:XT:36:LEU:CD1	20:XT:55:ILE:HD12	2.46	0.45
25:YA:70:G:H21	25:YA:71:A:N6	2.14	0.45
25:YA:1296:G:OP1	25:YA:2709:G:O2'	2.27	0.45
25:YA:2032:G:N2	28:YE:146:THR:HG23	2.28	0.45
25:YA:2208:U:O2'	27:YD:151:LYS:HG2	2.17	0.45
25:YA:2780:G:OP2	33:YN:118:LYS:HE2	2.17	0.45
25:YA:2875:C:H4'	39:YT:5:ALA:HB2	1.97	0.45
27:YD:109:ASP:HB2	27:YD:197:GLY:CA	2.46	0.45
28:YE:13:ARG:HH11	28:YE:13:ARG:HB3	1.82	0.45
28:YE:47:VAL:O	28:YE:47:VAL:HG23	2.16	0.45
35:YP:121:LYS:O	35:YP:123:LEU:N	2.49	0.45
38:YS:5:THR:OG1	38:YS:8:GLU:HG3	2.17	0.45
44:YY:101:LYS:HG2	44:YY:102:CYS:H	1.81	0.45
47:Y1:91:LYS:HA	47:Y1:91:LYS:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:Y2:65:ASN:O	48:Y2:66:GLU:C	2.59	0.45
1:QA:156:G:H1	1:QA:165:C:H42	1.64	0.45
1:QA:186:C:H5'	20:QT:78:ALA:HB1	1.97	0.45
1:QA:833:U:H2'	1:QA:834:C:H6	1.80	0.45
1:QA:1213:A:N1	1:QA:1215:G:H1'	2.31	0.45
1:QA:1330:U:H4'	13:QM:23:TYR:CE2	2.51	0.45
1:QA:1404:C:H2'	1:QA:1405:G:C8	2.52	0.45
7:QG:49:ILE:O	7:QG:53:LYS:HB3	2.16	0.45
10:QJ:55:LYS:CG	10:QJ:56:HIS:CD2	2.97	0.45
12:QL:17:LYS:HB2	12:QL:17:LYS:HZ2	1.81	0.45
25:RA:288:C:H2'	25:RA:289:A:H8	1.82	0.45
25:RA:675:A:N3	25:RA:2443:C:O2'	2.45	0.45
25:RA:2377:A:H2'	25:RA:2378:A:C8	2.51	0.45
25:RA:2635:C:H5"	28:RE:78:LEU:HA	1.99	0.45
27:RD:76:PRO:HB2	27:RD:116:GLN:OE1	2.17	0.45
27:RD:85:ASP:HA	27:RD:86:PRO:HD2	1.72	0.45
27:RD:118:VAL:HG22	27:RD:119:ALA:H	1.82	0.45
28:RE:22:PRO:CG	28:RE:22:PRO:O	2.63	0.45
31:RH:86:GLU:O	31:RH:87:LEU:CB	2.64	0.45
31:RH:94:TYR:CD1	31:RH:94:TYR:N	2.82	0.45
33:RN:19:GLU:HB2	33:RN:56:ASN:HD22	1.80	0.45
36:RQ:93:TYR:CD1	36:RQ:93:TYR:N	2.84	0.45
37:RR:103:ARG:NH1	42:RW:40:ASN:OD1	2.50	0.45
43:RX:67:GLY:C	43:RX:69:TYR:H	2.23	0.45
1:XA:323:U:O3'	20:XT:22:ARG:HD3	2.16	0.45
1:XA:537:G:H5"	12:XL:113:ARG:NH1	2.30	0.45
1:XA:877:C:H5"	8:XH:88:LYS:HD3	1.99	0.45
1:XA:923:A:H2'	1:XA:924:C:O4'	2.17	0.45
1:XA:954:G:H21	1:XA:1227:A:N6	1.91	0.45
1:XA:1127:G:H1'	1:XA:1147:C:H42	1.80	0.45
1:XA:1347:G:OP2	9:XI:107:ARG:HG2	2.16	0.45
3:XC:48:TYR:OH	3:XC:122:GLU:OE2	2.22	0.45
6:XF:100:ASN:O	18:XR:28:GLU:HG2	2.17	0.45
7:XG:45:ASP:O	7:XG:49:ILE:HG12	2.17	0.45
9:XI:114:TYR:N	9:XI:114:TYR:HD1	2.14	0.45
19:XS:47:HIS:O	19:XS:62:ILE:HG12	2.17	0.45
19:XS:63:THR:HG23	19:XS:66:MET:HG2	1.99	0.45
25:YA:270(T):G:OP1	47:Y1:97:LEU:HD13	2.16	0.45
25:YA:592:G:N3	54:Y8:4:MET:CE	2.79	0.45
25:YA:1754:C:P	39:YT:96:ARG:HH12	2.39	0.45
25:YA:2409:G:H2'	25:YA:2410:G:O4'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2563:U:H4'	34:YO:28:SER:HA	1.97	0.45
25:YA:2566:A:H4'	25:YA:2567:G:O5'	2.17	0.45
25:YA:2729:G:C2'	28:YE:187:ALA:HB2	2.46	0.45
27:YD:154:LYS:C	27:YD:155:LEU:HD12	2.41	0.45
28:YE:33:VAL:HG12	28:YE:90:THR:H	1.81	0.45
28:YE:95:ILE:HG22	28:YE:95:ILE:O	2.16	0.45
29:YF:196:LEU:O	29:YF:200:GLU:HG2	2.17	0.45
31:YH:66:GLY:O	31:YH:67:LEU:C	2.58	0.45
36:YQ:52:VAL:O	36:YQ:53:ALA:C	2.59	0.45
38:YS:78:LEU:HD21	38:YS:108:GLY:HA2	1.99	0.45
41:YV:19:LYS:HG3	41:YV:95:LEU:HD23	1.98	0.45
41:YV:36:PRO:HA	41:YV:56:SER:OG	2.16	0.45
44:YY:89:PHE:C	44:YY:90:LEU:HD13	2.41	0.45
45:YZ:128:VAL:HG22	45:YZ:129:SER:H	1.82	0.45
45:YZ:141:VAL:HG23	45:YZ:144:LEU:HB2	1.98	0.45
1:QA:109:A:H2'	1:QA:326:G:N2	2.32	0.45
1:QA:636:U:H5'	17:QQ:2:PRO:HG3	1.98	0.45
1:QA:1304:G:N1	1:QA:1332:A:OP2	2.42	0.45
1:QA:1348:U:C4	1:QA:1374:A:H2	2.33	0.45
7:QG:113:GLU:CG	7:QG:119:ARG:HG2	2.47	0.45
9:QI:126:SER:O	9:QI:128:ARG:N	2.45	0.45
11:QK:91:ARG:NH1	11:QK:110:ASP:OD2	2.48	0.45
12:QL:6:THR:H	12:QL:9:GLN:NE2	1.97	0.45
20:QT:33:ILE:HD12	20:QT:63:ILE:HG12	1.96	0.45
25:RA:78:A:H2'	25:RA:79:G:C8	2.52	0.45
25:RA:856:C:H42	25:RA:921:G:H1	1.63	0.45
25:RA:1186:G:H2'	25:RA:1187:G:O4'	2.16	0.45
25:RA:2364:C:OP1	46:R0:55:ARG:NH1	2.50	0.45
25:RA:2568:C:H2'	25:RA:2569:G:O4'	2.16	0.45
25:RA:2593:U:H2'	25:RA:2594:C:C6	2.51	0.45
27:RD:105:ILE:HD12	27:RD:105:ILE:HA	1.63	0.45
30:RG:10:LYS:O	30:RG:14:GLU:HB3	2.17	0.45
30:RG:102:PHE:O	30:RG:106:LEU:N	2.50	0.45
31:RH:6:ARG:C	31:RH:8:PRO:CD	2.90	0.45
32:RI:14:ASP:O	32:RI:16:GLY:N	2.50	0.45
32:RI:29:TYR:C	32:RI:32:PRO:HD2	2.41	0.45
34:RO:48:PRO:O	34:RO:49:ARG:HG2	2.17	0.45
38:RS:83:LYS:C	38:RS:109:GLY:HA3	2.42	0.45
45:RZ:108:PRO:HB2	45:RZ:109:ALA:H	1.62	0.45
48:R2:47:ASN:O	48:R2:48:HIS:CG	2.70	0.45
1:XA:767:A:O2'	1:XA:1524:C:O2	2.33	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:971:G:N2	1:XA:1363:A:OP2	2.33	0.45
1:XA:1410:G:H2'	1:XA:1411:C:H6	1.82	0.45
2:XB:120:ALA:C	2:XB:122:PHE:H	2.25	0.45
15:XO:32:LEU:HD11	15:XO:62:GLN:HG2	1.99	0.45
16:XP:17:TYR:CE2	16:XP:41:PRO:HG3	2.52	0.45
25:YA:855:G:O2'	46:Y0:27:GLU:OE2	2.24	0.45
25:YA:1869:G:H5'	25:YA:1870:C:OP2	2.16	0.45
25:YA:1930:G:HO2'	25:YA:1931:U:P	2.40	0.45
25:YA:2756:U:H4'	25:YA:2757:A:OP1	2.16	0.45
27:YD:198:ASN:ND2	27:YD:198:ASN:O	2.50	0.45
28:YE:199:ARG:HH11	28:YE:199:ARG:HG3	1.82	0.45
29:YF:155:LEU:HA	29:YF:174:VAL:CG1	2.46	0.45
33:YN:8:GLN:C	33:YN:9:VAL:CG1	2.86	0.45
36:YQ:26:TYR:O	36:YQ:27:VAL:O	2.34	0.45
44:YY:94:LYS:HD2	44:YY:101:LYS:HZ3	1.82	0.45
54:Y8:52:LYS:O	54:Y8:52:LYS:CG	2.64	0.45
1:QA:607:A:C2	16:QP:31:LYS:HB2	2.52	0.45
1:QA:745:C:OP1	1:QA:851:G:O2'	2.35	0.45
1:QA:804:U:H5"	1:QA:805:C:OP2	2.17	0.45
1:QA:920:U:H2'	1:QA:921:U:H6	1.81	0.45
1:QA:1032(A):G:H2'	1:QA:1032(B):G:C8	2.52	0.45
1:QA:1126:U:OP2	1:QA:1281:U:H1'	2.16	0.45
2:QB:30:ARG:HH21	2:QB:194:PRO:HG2	1.81	0.45
2:QB:76:GLN:O	2:QB:208:ILE:HG12	2.17	0.45
8:QH:104:ARG:O	8:QH:107:LEU:HB2	2.16	0.45
25:RA:482:A:H4'	44:RY:47:LYS:HD2	1.99	0.45
25:RA:919:G:N2	25:RA:2269:A:OP2	2.50	0.45
25:RA:1085:A:H2'	25:RA:1086:A:C4	2.52	0.45
25:RA:1444(A):A:H4'	25:RA:1460:A:O2'	2.17	0.45
25:RA:2795:G:H3'	25:RA:2797:U:C5'	2.47	0.45
26:RB:38:C:N3	26:RB:44:G:N2	2.58	0.45
37:RR:42:LYS:HA	37:RR:45:ARG:HD2	1.98	0.45
37:RR:78:LYS:HE2	37:RR:83:ILE:HD11	1.98	0.45
51:R5:15:ARG:HA	51:R5:18:ALA:HB3	1.99	0.45
54:R8:48:PHE:N	54:R8:48:PHE:HD1	2.14	0.45
1:XA:781:A:O2'	1:XA:1522:U:O2	2.32	0.45
1:XA:1314:C:OP1	19:XS:6:LYS:HE3	2.17	0.45
2:XB:115:LEU:HD13	2:XB:145:LEU:HB3	1.98	0.45
12:XL:126:LYS:HB2	12:XL:126:LYS:HZ3	1.81	0.45
25:YA:312:G:H5'	25:YA:331:A:O2'	2.17	0.45
25:YA:888:C:H3'	25:YA:889:C:C4'	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:1665:A:H1'	34:YO:1:MET:HG3	1.98	0.45
25:YA:1717:G:H1	25:YA:1742:C:H42	1.65	0.45
26:YB:8:U:O2'	38:YS:40:ILE:CD1	2.65	0.45
26:YB:43:C:H5'	50:Y4:1:MET:HA	1.97	0.45
27:YD:25:THR:CG2	27:YD:25:THR:O	2.65	0.45
27:YD:35:LYS:CE	27:YD:64:ILE:C	2.89	0.45
27:YD:68:LYS:HD2	27:YD:70:TRP:CZ2	2.52	0.45
27:YD:92:ILE:HD12	27:YD:104:TYR:HD2	1.82	0.45
28:YE:2:LYS:O	28:YE:199:ARG:HA	2.17	0.45
29:YF:198:ALA:C	29:YF:200:GLU:H	2.25	0.45
35:YP:65:ARG:HB2	54:Y8:12:LYS:O	2.17	0.45
39:YT:102:ILE:HA	39:YT:105:LEU:CD2	2.47	0.45
40:YU:90:VAL:HG13	41:YV:4:ILE:HG21	1.98	0.45
42:YW:97:LYS:HE2	42:YW:99:ARG:NH2	2.31	0.45
46:Y0:23:VAL:HA	46:Y0:38:VAL:HA	1.99	0.45
47:Y1:79:GLY:N	47:Y1:80:LEU:HD23	2.32	0.45
52:Y6:28:ARG:HB3	52:Y6:30:THR:C	2.41	0.45
1:QA:73:G:H2'	1:QA:74:C:O4'	2.16	0.45
1:QA:458:C:H42	1:QA:474:G:H1	1.63	0.45
2:QB:163:PHE:HA	2:QB:185:ILE:HG13	1.98	0.45
2:QB:178:ARG:HH21	8:QH:74:PRO:HG3	1.82	0.45
5:QE:151:LEU:HD22	8:QH:67:PRO:HD3	1.99	0.45
6:QF:45:LEU:HD12	6:QF:59:TYR:HD2	1.82	0.45
10:QJ:40:LEU:HB2	10:QJ:69:ASN:CB	2.47	0.45
25:RA:198:C:O2'	25:RA:199:A:H5'	2.15	0.45
25:RA:1024:G:C6	25:RA:1025:G:C6	3.05	0.45
25:RA:1111:A:H4'	31:RH:3:ARG:HD2	1.98	0.45
25:RA:1188:U:O2'	25:RA:1189:A:H5'	2.15	0.45
25:RA:2294:C:OP2	38:RS:13:ARG:NH1	2.50	0.45
28:RE:2:LYS:HG2	28:RE:95:ILE:HG22	1.99	0.45
28:RE:2:LYS:O	28:RE:199:ARG:HA	2.17	0.45
34:RO:22:ILE:HG12	34:RO:41:ALA:HA	1.98	0.45
36:RQ:119:ARG:HH11	36:RQ:119:ARG:CG	2.25	0.45
38:RS:10:ARG:O	38:RS:14:VAL:HG12	2.17	0.45
1:XA:358:U:H2'	1:XA:359:U:O4'	2.17	0.45
1:XA:376:G:H5"	16:XP:5:ARG:HD2	1.99	0.45
3:XC:153:VAL:HG22	3:XC:198:VAL:HG22	1.98	0.45
4:XD:86:LYS:HD2	4:XD:86:LYS:H	1.82	0.45
10:XJ:16:LEU:HD11	10:XJ:70:ARG:HB2	1.99	0.45
10:XJ:55:LYS:HG3	10:XJ:56:HIS:CG	2.51	0.45
12:XL:117:ARG:NH2	12:XL:124:LYS:HD3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:XM:65:LYS:O	13:XM:70:LEU:HD23	2.17	0.45
19:XS:66:MET:HB2	19:XS:74:PHE:CZ	2.52	0.45
25:YA:192:C:H2'	25:YA:193:U:H5'	1.98	0.45
25:YA:612:G:O2'	25:YA:616:A:N1	2.45	0.45
25:YA:754:C:H2'	25:YA:755:C:H6	1.82	0.45
25:YA:828:U:H4'	25:YA:831:G:N1	2.31	0.45
25:YA:844:C:H2'	25:YA:845:G:O4'	2.17	0.45
25:YA:984:A:H5"	25:YA:985:C:H5	1.82	0.45
25:YA:1130:U:O2'	25:YA:1131:G:O5'	2.30	0.45
25:YA:1332:G:N2	25:YA:1609:A:HO2'	2.15	0.45
25:YA:2344:U:OP1	52:Y6:38:LYS:HD3	2.17	0.45
26:YB:77:U:OP1	45:YZ:19:ARG:NH2	2.49	0.45
27:YD:226:MET:H	27:YD:226:MET:HG2	1.53	0.45
28:YE:4:ILE:HG12	28:YE:91:VAL:HG11	1.99	0.45
28:YE:21:VAL:HG23	28:YE:22:PRO:CD	2.46	0.45
28:YE:154:LYS:C	28:YE:154:LYS:HD3	2.42	0.45
29:YF:61:GLY:O	29:YF:62:ARG:C	2.57	0.45
36:YQ:5:ARG:O	36:YQ:6:ARG:O	2.35	0.45
36:YQ:30:GLY:CA	36:YQ:107:ALA:HB2	2.39	0.45
36:YQ:34:LEU:HB2	36:YQ:118:LEU:HD22	1.99	0.45
38:YS:111:GLU:O	38:YS:112:PHE:CD2	2.70	0.45
39:YT:6:LEU:HA	39:YT:9:LEU:HB2	1.99	0.45
48:Y2:28:LYS:HB3	48:Y2:57:ILE:HG12	1.98	0.45
1:QA:6:G:N2	5:QE:98:THR:OG1	2.50	0.45
1:QA:1162:C:N4	1:QA:1174:G:H1	2.12	0.45
1:QA:1299:A:C8	1:QA:1301:U:H1'	2.52	0.45
1:QA:1465:C:H2'	1:QA:1466:C:O4'	2.17	0.45
8:QH:6:ILE:O	8:QH:10:LEU:HG	2.17	0.45
15:QO:4:THR:HB	15:QO:6:GLU:CD	2.41	0.45
25:RA:67:U:O4	25:RA:74:A:N1	2.49	0.45
25:RA:443:A:H3'	29:RF:45:ARG:NH1	2.31	0.45
25:RA:2443:C:H2'	25:RA:2444:G:H8	1.82	0.45
25:RA:2452:C:O4'	56:Z5:101:PPU:HB2	2.17	0.45
25:RA:2527:C:H5"	55:R9:30:PRO:HB2	1.99	0.45
25:RA:2660:A:H2'	25:RA:2661:G:O4'	2.17	0.45
25:RA:2760:C:H2'	25:RA:2761:G:H5"	1.98	0.45
28:RE:93:VAL:C	28:RE:95:ILE:N	2.75	0.45
35:RP:61:ARG:HG3	54:R8:13:ARG:HD3	1.97	0.45
47:R1:83:GLU:OE2	47:R1:83:GLU:N	2.49	0.45
52:R6:34:LEU:HD13	52:R6:34:LEU:H	1.82	0.45
54:R8:36:LYS:HB3	54:R8:40:GLU:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:107:G:C2	1:XA:108:G:H1'	2.52	0.45
1:XA:560:U:O2'	1:XA:561:U:OP2	2.33	0.45
1:XA:729:A:H2'	1:XA:730:G:C8	2.51	0.45
1:XA:1247:U:H1'	1:XA:1291:G:N2	2.31	0.45
8:XH:51:VAL:HG11	8:XH:60:ARG:HG3	1.99	0.45
11:XK:28:THR:OG1	11:XK:90:GLY:HA3	2.17	0.45
22:XV:51:C:H2'	22:XV:52:G:O4'	2.17	0.45
25:YA:442:G:C4'	29:YF:46:ARG:HD3	2.47	0.45
25:YA:570:G:H2'	25:YA:2030:A:C6	2.52	0.45
25:YA:669:G:H2'	25:YA:669:G:N3	2.32	0.45
25:YA:1162:G:O2'	41:YV:90:PRO:HG2	2.16	0.45
25:YA:1188:U:H4'	41:YV:79:VAL:HG22	1.99	0.45
25:YA:1510:A:H2'	25:YA:1510:A:N3	2.32	0.45
25:YA:2578:G:N7	28:YE:140:SER:HB2	2.31	0.45
26:YB:30:C:OP2	38:YS:32:LEU:HD11	2.17	0.45
36:YQ:58:PHE:O	36:YQ:58:PHE:CD1	2.70	0.45
36:YQ:93:TYR:CD1	36:YQ:93:TYR:N	2.85	0.45
50:Y4:16:CYS:HB3	50:Y4:33:VAL:HB	1.98	0.45
54:Y8:36:LYS:HB3	54:Y8:40:GLU:HG2	1.99	0.45
1:QA:756:C:H2'	1:QA:757:U:O4'	2.17	0.45
1:QA:900:A:H2'	1:QA:901:A:C8	2.52	0.45
8:QH:6:ILE:HB	8:QH:85:ARG:HH11	1.82	0.45
11:QK:41:THR:HG21	11:QK:71:LYS:HB2	1.99	0.45
25:RA:300:A:H2'	25:RA:334:C:H1'	1.99	0.45
25:RA:1353:A:O3'	27:RD:38:LYS:NZ	2.36	0.45
25:RA:1728:G:H3'	25:RA:1729:A:C5'	2.46	0.45
25:RA:1819:A:H2'	27:RD:178:PRO:HB2	1.98	0.45
25:RA:2350:C:H2'	25:RA:2351:G:O4'	2.17	0.45
28:RE:77:ILE:O	28:RE:78:LEU:O	2.34	0.45
34:RO:87:ILE:HD12	34:RO:91:LEU:HD12	1.99	0.45
34:RO:111:PHE:HB3	34:RO:114:ILE:HG13	1.98	0.45
36:RQ:10:ARG:O	36:RQ:11:LYS:CB	2.64	0.45
36:RQ:11:LYS:HE2	36:RQ:87:LYS:HA	1.98	0.45
43:RX:15:GLU:H	43:RX:15:GLU:CD	2.25	0.45
44:RY:68:HIS:CE1	44:RY:70:SER:HB3	2.52	0.45
48:R2:41:ILE:HD11	48:R2:44:LEU:HB2	1.99	0.45
54:R8:9:GLY:O	54:R8:13:ARG:HG2	2.16	0.45
1:XA:322:C:O2'	20:XT:23:ARG:HD2	2.17	0.45
1:XA:707:C:OP1	11:XK:85:ARG:NH1	2.49	0.45
1:XA:1436:U:OP1	20:XT:23:ARG:NH2	2.44	0.45
4:XD:25:ARG:NH1	4:XD:30:LYS:HG3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:XI:111:ARG:HH22	10:XJ:62:HIS:CE1	2.35	0.45
19:XS:24:ALA:O	19:XS:25:LYS:HB3	2.17	0.45
25:YA:1778:U:H2'	25:YA:1784:A:N6	2.32	0.45
25:YA:2212:A:H1'	25:YA:2215:G:C4	2.52	0.45
25:YA:2572:A:C8	28:YE:144:ARG:HB3	2.52	0.45
25:YA:2671:A:H2'	25:YA:2672:G:O4'	2.16	0.45
27:YD:214:TRP:C	27:YD:216:GLY:H	2.25	0.45
29:YF:144:LYS:C	29:YF:146:ALA:N	2.75	0.45
31:YH:6:ARG:C	31:YH:8:PRO:CD	2.89	0.45
31:YH:137:ASP:HB2	31:YH:140:LYS:HE3	1.98	0.45
38:YS:3:ARG:O	38:YS:4:LEU:O	2.35	0.45
50:Y4:22:ILE:HG22	50:Y4:23:GLU:H	1.82	0.45
1:QA:474:G:H2'	1:QA:475:G:C8	2.52	0.44
1:QA:696:A:N1	1:QA:797:C:O2'	2.41	0.44
1:QA:1072:G:H2'	1:QA:1073:U:C6	2.52	0.44
1:QA:1210:C:H3'	1:QA:1211:U:H5"	1.98	0.44
6:QF:41:GLU:HB3	6:QF:62:TRP:HB3	1.99	0.44
12:QL:15:ARG:HG2	17:QQ:32:TYR:OH	2.17	0.44
12:QL:126:LYS:HB2	12:QL:126:LYS:HZ2	1.82	0.44
25:RA:321:G:OP1	29:RF:135:LYS:NZ	2.48	0.44
25:RA:395:U:H2'	25:RA:396:G:C8	2.52	0.44
25:RA:1048:A:H2	25:RA:1112:G:H21	1.64	0.44
25:RA:2451:A:N1	56:Z5:101:PPU:HE2	2.31	0.44
28:RE:4:ILE:HG12	28:RE:91:VAL:HG11	1.99	0.44
28:RE:36:ARG:HH11	28:RE:36:ARG:CB	2.28	0.44
28:RE:154:LYS:HD3	28:RE:154:LYS:C	2.42	0.44
29:RF:149:ASP:OD1	29:RF:149:ASP:N	2.45	0.44
31:RH:149:ARG:HA	31:RH:162:ILE:HG21	1.99	0.44
37:RR:28:LEU:HD12	37:RR:48:VAL:HG11	1.99	0.44
37:RR:29:LEU:HD12	37:RR:29:LEU:HA	1.74	0.44
50:R4:6:HIS:O	50:R4:7:PRO:C	2.61	0.44
50:R4:33:VAL:CG1	50:R4:34:GLU:H	2.22	0.44
54:R8:16:ILE:CD1	54:R8:57:ARG:HG2	2.42	0.44
54:R8:40:GLU:C	54:R8:42:ARG:N	2.68	0.44
1:XA:255:G:H2'	1:XA:256:U:C6	2.52	0.44
1:XA:347:G:O2'	1:XA:348:G:OP2	2.29	0.44
1:XA:792:A:H4'	1:XA:793:U:O5'	2.18	0.44
1:XA:1336:C:HO2'	1:XA:1337:G:P	2.40	0.44
8:XH:75:ARG:HA	8:XH:76:PRO:HD2	1.71	0.44
13:XM:121:LYS:HE2	13:XM:121:LYS:HA	1.98	0.44
25:YA:1693:U:H1'	27:YD:14:ARG:HH22	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2584:U:H5'	56:Z6:101:PPU:C10	2.45	0.44
26:YB:81:G:O6	26:YB:95:U:C2	2.70	0.44
27:YD:52:ARG:HB2	27:YD:53:PHE:CD2	2.52	0.44
32:YI:93:THR:O	32:YI:97:ILE:HG12	2.17	0.44
34:YO:17:ARG:NH2	34:YO:47:ILE:HD13	2.32	0.44
34:YO:68:GLU:H	34:YO:68:GLU:CD	2.24	0.44
36:YQ:65:PHE:O	36:YQ:66:ILE:CG1	2.48	0.44
36:YQ:133:ARG:CG	36:YQ:134:ARG:N	2.78	0.44
38:YS:78:LEU:HD21	38:YS:108:GLY:CA	2.47	0.44
38:YS:83:LYS:CE	38:YS:109:GLY:HA2	2.47	0.44
50:Y4:43:TYR:C	50:Y4:43:TYR:HD1	2.25	0.44
1:QA:333:G:H4'	20:QT:16:HIS:NE2	2.32	0.44
1:QA:345:C:H5'	1:QA:346:G:C5	2.53	0.44
1:QA:437:U:H2'	1:QA:438:G:O4'	2.16	0.44
1:QA:636:U:H2'	1:QA:637:G:C8	2.52	0.44
1:QA:895:G:H2'	1:QA:896:C:C6	2.52	0.44
1:QA:1436:U:H2'	1:QA:1437:C:O4'	2.17	0.44
3:QC:34:LEU:HG	14:QN:25:VAL:HG11	2.00	0.44
4:QD:175:SER:HB3	4:QD:186:LEU:HD11	1.99	0.44
5:QE:103:GLY:O	5:QE:107:ARG:HB3	2.17	0.44
12:QL:120:TYR:N	12:QL:120:TYR:CD1	2.85	0.44
19:QS:10:PHE:HB2	19:QS:39:THR:H	1.82	0.44
25:RA:135:G:H1	25:RA:144:C:H42	1.66	0.44
25:RA:859:G:O2'	25:RA:860:U:OP2	2.35	0.44
25:RA:1510:A:N3	25:RA:1510:A:H2'	2.32	0.44
27:RD:70:TRP:HZ3	27:RD:146:GLU:OE2	2.01	0.44
28:RE:36:ARG:O	28:RE:37:ARG:C	2.59	0.44
31:RH:51:ARG:HG3	31:RH:51:ARG:NH1	2.30	0.44
33:RN:114:ARG:O	33:RN:115:ARG:HB3	2.17	0.44
40:RU:65:ILE:HG12	40:RU:96:ALA:CB	2.47	0.44
45:RZ:58:VAL:O	45:RZ:60:GLU:N	2.47	0.44
51:R5:56:LYS:O	51:R5:58:LEU:N	2.50	0.44
54:R8:17:THR:O	54:R8:20:GLY:N	2.46	0.44
1:XA:983:A:H5"	1:XA:984:C:OP2	2.17	0.44
1:XA:1288:A:N3	1:XA:1352:C:O2'	2.43	0.44
5:XE:110:LEU:HD13	5:XE:118:ILE:HD13	1.98	0.44
7:XG:138:LYS:HE2	7:XG:142:GLU:OE2	2.17	0.44
12:XL:120:TYR:N	12:XL:120:TYR:CD1	2.86	0.44
19:XS:81:ARG:HE	19:XS:81:ARG:HB2	1.36	0.44
25:YA:658:C:H2'	25:YA:659:C:C6	2.52	0.44
27:YD:80:ALA:O	27:YD:113:VAL:HG13	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:166:GLN:HA	27:YD:166:GLN:NE2	2.32	0.44
27:YD:177:LEU:O	27:YD:179:SER:N	2.51	0.44
28:YE:121:ASN:O	28:YE:122:PHE:C	2.57	0.44
29:YF:149:ASP:OD2	29:YF:151:SER:HB3	2.17	0.44
42:YW:33:ARG:NH2	42:YW:52:GLU:OE1	2.50	0.44
51:Y5:40:LYS:HZ3	51:Y5:46:CYS:C	2.26	0.44
1:QA:67:C:O2'	1:QA:171:A:N3	2.40	0.44
1:QA:1327:C:H2'	1:QA:1328:C:C6	2.52	0.44
2:QB:51:LEU:HD22	2:QB:55:PHE:HE2	1.82	0.44
10:QJ:33:GLN:O	10:QJ:75:ILE:HG12	2.17	0.44
25:RA:847:U:O4	25:RA:933:A:C6	2.70	0.44
25:RA:1266:G:OP2	51:R5:20:ARG:NE	2.44	0.44
28:RE:101:ARG:HD2	28:RE:171:GLU:HA	1.98	0.44
29:RF:9:ILE:HG23	29:RF:20:LEU:O	2.18	0.44
30:RG:6:ALA:HB2	50:R4:23:GLU:CD	2.41	0.44
30:RG:37:VAL:O	30:RG:94:LEU:HG	2.17	0.44
31:RH:53:GLU:OE1	31:RH:53:GLU:HA	2.16	0.44
33:RN:116:LEU:HD23	33:RN:116:LEU:HA	1.78	0.44
44:RY:42:VAL:HG12	44:RY:65:ALA:HB3	1.99	0.44
50:R4:51:ASP:O	50:R4:51:ASP:CG	2.60	0.44
52:R6:32:ASN:N	52:R6:32:ASN:OD1	2.49	0.44
1:XA:343:U:O2	1:XA:346:G:N1	2.47	0.44
1:XA:575:G:O2'	1:XA:821:G:OP2	2.33	0.44
1:XA:1162:C:N3	1:XA:1174:G:N2	2.57	0.44
1:XA:1285:A:H4'	1:XA:1286:A:O5'	2.17	0.44
1:XA:1399:C:C2	1:XA:1502:A:N6	2.85	0.44
4:XD:26:CYS:SG	4:XD:31:CYS:CB	3.02	0.44
9:XI:25:LYS:HE3	9:XI:25:LYS:HB2	1.72	0.44
9:XI:95:LYS:HZ3	9:XI:96:LEU:HD13	1.83	0.44
21:XU:5:ASP:HB3	21:XU:8:THR:OG1	2.17	0.44
25:YA:607:U:OP1	29:YF:102:PRO:HA	2.17	0.44
25:YA:1407:C:N4	25:YA:1595:G:H1	2.12	0.44
25:YA:1469:A:H2'	25:YA:1470:G:O4'	2.16	0.44
25:YA:1491:G:O2'	27:YD:101:GLU:HB2	2.18	0.44
25:YA:1803:A:C3'	27:YD:259:THR:HG21	2.47	0.44
27:YD:166:GLN:CA	27:YD:166:GLN:NE2	2.78	0.44
28:YE:93:VAL:C	28:YE:95:ILE:N	2.75	0.44
29:YF:33:LEU:HD23	35:YP:1:MET:SD	2.57	0.44
34:YO:86:ILE:HG22	34:YO:94:ARG:HD3	2.00	0.44
36:YQ:66:ILE:O	36:YQ:104:PHE:N	2.49	0.44
38:YS:56:LEU:O	38:YS:57:LYS:O	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:YT:42:ILE:HG21	39:YT:84:GLN:NE2	2.32	0.44
39:YT:61:PHE:CE1	39:YT:76:PHE:HB2	2.53	0.44
40:YU:104:GLN:OE1	40:YU:105:VAL:HG23	2.17	0.44
43:YX:35:THR:O	43:YX:39:ILE:HG13	2.16	0.44
54:Y8:47:LYS:HD2	54:Y8:48:PHE:N	2.33	0.44
1:QA:765:G:N2	1:QA:813:U:OP2	2.45	0.44
1:QA:768:A:N3	1:QA:1512:U:O2'	2.46	0.44
12:QL:120:TYR:O	12:QL:121:GLY:O	2.36	0.44
19:QS:5:LEU:HD11	50:R4:66:SER:HB2	1.99	0.44
19:QS:66:MET:HB2	19:QS:74:PHE:CZ	2.52	0.44
25:RA:122:G:OP1	25:RA:149:A:O2'	2.29	0.44
25:RA:1266:G:O2'	25:RA:2012:G:O6	2.29	0.44
25:RA:2031:A:N3	25:RA:2455:G:O2'	2.39	0.44
25:RA:2567:G:H2'	25:RA:2568:C:C6	2.53	0.44
25:RA:2774:C:H2'	25:RA:2775:A:O4'	2.17	0.44
29:RF:127:GLU:C	29:RF:129:PHE:H	2.26	0.44
32:RI:5:LEU:HD23	32:RI:9:LEU:HD11	1.99	0.44
35:RP:52:GLU:OE1	35:RP:55:ARG:NH1	2.50	0.44
38:RS:11:LYS:HG3	38:RS:91:PRO:HD3	1.98	0.44
50:R4:15:ILE:CG2	50:R4:20:ASN:ND2	2.81	0.44
54:R8:47:LYS:HD2	54:R8:48:PHE:N	2.33	0.44
1:XA:118:U:H3'	1:XA:288:A:H61	1.82	0.44
1:XA:302:G:N3	1:XA:556:C:H4'	2.33	0.44
1:XA:564:C:H5'	17:XQ:32:TYR:CE1	2.52	0.44
1:XA:738:C:H2'	1:XA:739:C:H6	1.82	0.44
7:XG:87:VAL:HG11	7:XG:155:ARG:HA	1.99	0.44
8:XH:65:TYR:HA	8:XH:79:VAL:HG23	1.98	0.44
8:XH:83:ILE:HB	8:XH:137:VAL:HG13	1.99	0.44
20:XT:87:LYS:O	20:XT:91:LEU:HG	2.18	0.44
22:XV:3:G:O2'	22:XV:4:G:H8	1.99	0.44
22:XV:36:U:H2'	22:XV:37:A:C8	2.53	0.44
25:YA:242:G:C8	54:Y8:5:LYS:HG2	2.52	0.44
25:YA:242:G:H1'	25:YA:243:U:OP2	2.18	0.44
25:YA:264:C:H2'	25:YA:265:A:H5"	1.99	0.44
25:YA:459:U:H2'	25:YA:460:A:H8	1.83	0.44
25:YA:1889:A:N1	25:YA:2234:G:H1'	2.32	0.44
25:YA:2572:A:C2	28:YE:144:ARG:NH2	2.85	0.44
28:YE:25:VAL:HG11	39:YT:11:GLU:CD	2.43	0.44
28:YE:50:GLY:CA	28:YE:74:PRO:HG3	2.46	0.44
28:YE:172:VAL:HG13	28:YE:182:LEU:HD11	1.98	0.44
29:YF:111:ALA:O	29:YF:112:MET:C	2.59	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:YU:66:ASN:HB2	40:YU:76:TYR:HB2	1.99	0.44
47:Y1:53:VAL:HB	47:Y1:58:ILE:HD13	1.99	0.44
48:Y2:41:ILE:HD12	48:Y2:41:ILE:O	2.16	0.44
52:Y6:7:ILE:HD12	52:Y6:7:ILE:HA	1.85	0.44
52:Y6:34:LEU:HD13	52:Y6:34:LEU:H	1.82	0.44
1:QA:465:A:N6	1:QA:467:G:C6	2.86	0.44
1:QA:483:C:OP2	1:QA:484:G:O2'	2.24	0.44
1:QA:514:C:H2'	1:QA:515:G:H8	1.82	0.44
1:QA:555:C:H2'	1:QA:556:C:C6	2.53	0.44
1:QA:665:A:H2'	1:QA:725:G:N2	2.32	0.44
1:QA:860:A:H2'	1:QA:861:G:O4'	2.17	0.44
6:QF:23:LYS:O	6:QF:27:GLN:HG2	2.17	0.44
10:QJ:55:LYS:CE	10:QJ:56:HIS:NE2	2.73	0.44
12:QL:91:LYS:HE2	12:QL:91:LYS:HB2	1.76	0.44
15:QO:39:LEU:HA	15:QO:39:LEU:HD23	1.69	0.44
25:RA:508:G:HO2'	25:RA:509:C:P	2.41	0.44
25:RA:573:G:O2'	25:RA:574:C:H3'	2.18	0.44
25:RA:1248:G:C5	40:RU:3:ARG:HB2	2.52	0.44
25:RA:1286:A:HO2'	25:RA:1288:U:P	2.39	0.44
25:RA:1542:G:H5"	25:RA:1543:A:OP2	2.18	0.44
25:RA:1628:G:H1	25:RA:1638:C:N4	2.14	0.44
25:RA:2688:U:H2'	25:RA:2719:G:N2	2.32	0.44
26:RB:52:A:H62	38:RS:33:LYS:HG3	1.82	0.44
26:RB:83:G:H4'	49:R3:52:HIS:CG	2.53	0.44
28:RE:143:ASN:ND2	28:RE:143:ASN:N	2.65	0.44
30:RG:51:ARG:O	30:RG:53:LEU:N	2.48	0.44
30:RG:95:ARG:O	30:RG:99:MET:HG2	2.17	0.44
31:RH:84:SER:OG	31:RH:85:LYS:N	2.51	0.44
1:XA:738:C:H2'	1:XA:739:C:C6	2.52	0.44
1:XA:1236:A:H4'	1:XA:1304:G:H4'	1.99	0.44
2:XB:33:TYR:HB2	2:XB:43:ASP:HB2	1.99	0.44
5:XE:9:LYS:HB2	5:XE:9:LYS:HE3	1.89	0.44
5:XE:41:VAL:HG22	5:XE:113:ALA:HB2	1.99	0.44
10:XJ:61:GLU:OE1	14:XN:58:LYS:HE2	2.17	0.44
13:XM:115:LYS:HE3	13:XM:115:LYS:HB2	1.75	0.44
18:XR:52:PRO:HB2	18:XR:54:ARG:HG2	2.00	0.44
25:YA:593:G:C1'	54:Y8:4:MET:HE1	2.44	0.44
25:YA:769:G:H5'	25:YA:1379:A:N6	2.33	0.44
25:YA:1035:U:H2'	25:YA:1036:G:H8	1.82	0.44
25:YA:1899:G:H21	25:YA:1902:C:H42	1.65	0.44
25:YA:2395:C:O2'	47:Y1:30:VAL:HG12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:2469:A:H2	25:YA:2481:G:H21	1.65	0.44
25:YA:2596:U:H2'	25:YA:2597:G:O4'	2.18	0.44
27:YD:213:ARG:HD2	27:YD:213:ARG:HA	1.60	0.44
28:YE:1:MET:HE2	28:YE:1:MET:H1	1.83	0.44
29:YF:24:LEU:HD12	29:YF:24:LEU:N	2.33	0.44
29:YF:65:TRP:HZ3	29:YF:73:ALA:C	2.26	0.44
29:YF:201:VAL:HG13	29:YF:202:PHE:N	2.33	0.44
30:YG:114:ILE:HB	30:YG:117:PHE:HB2	2.00	0.44
33:YN:134:ARG:O	33:YN:136:GLU:N	2.50	0.44
35:YP:65:ARG:HE	54:Y8:15:LYS:HB2	1.81	0.44
41:YV:99:ILE:HD13	41:YV:99:ILE:H	1.82	0.44
45:YZ:98:MET:HE3	45:YZ:98:MET:HB2	1.92	0.44
2:QB:71:VAL:HA	2:QB:93:VAL:HB	2.00	0.44
2:QB:228:GLY:O	2:QB:230:VAL:N	2.50	0.44
3:QC:70:VAL:HG21	3:QC:76:VAL:HG11	2.00	0.44
11:QK:92:GLU:HB3	11:QK:96:ARG:NH1	2.33	0.44
11:QK:96:ARG:HA	11:QK:99:GLN:HE21	1.83	0.44
12:QL:117:ARG:NH2	12:QL:124:LYS:HD3	2.32	0.44
15:QO:50:HIS:O	15:QO:53:HIS:HB3	2.17	0.44
20:QT:36:LEU:CD1	20:QT:55:ILE:HG23	2.48	0.44
25:RA:1336:A:H2'	25:RA:1337:G:C8	2.53	0.44
25:RA:1592:C:H2'	25:RA:1593:G:H8	1.83	0.44
27:RD:49:ILE:CD1	27:RD:52:ARG:HA	2.47	0.44
28:RE:95:ILE:HG22	28:RE:95:ILE:O	2.16	0.44
34:RO:106:LEU:HA	34:RO:106:LEU:HD23	1.81	0.44
41:RV:35:LEU:CD2	41:RV:57:VAL:HG22	2.47	0.44
41:RV:49:THR:HB	41:RV:50:PRO:HD2	1.99	0.44
43:RX:57:LEU:HD11	43:RX:78:LYS:HD2	1.99	0.44
47:R1:49:VAL:HG11	47:R1:70:VAL:HG11	1.98	0.44
50:R4:68:ARG:HH11	50:R4:69:LYS:HG2	1.82	0.44
55:R9:24:TYR:CE2	55:R9:35:ARG:HG3	2.53	0.44
1:XA:21:G:H2'	1:XA:22:G:C8	2.53	0.44
1:XA:260:G:H2'	1:XA:261:U:C6	2.53	0.44
1:XA:509:A:H2'	1:XA:510:A:C8	2.53	0.44
1:XA:1069:C:H42	1:XA:1106:G:H1	1.66	0.44
1:XA:1128:C:C2	1:XA:1144:G:N2	2.86	0.44
1:XA:1262:C:H2'	1:XA:1263:C:C6	2.52	0.44
1:XA:1320:C:C4	19:XS:36:ARG:HG3	2.53	0.44
15:XO:82:ILE:O	15:XO:86:GLY:N	2.51	0.44
16:XP:18:ARG:NH1	16:XP:32:TYR:OH	2.50	0.44
20:XT:44:ALA:HB1	20:XT:91:LEU:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:102:G:OP2	48:Y2:7:ARG:NH2	2.51	0.44
25:YA:459:U:H2'	25:YA:460:A:C8	2.52	0.44
25:YA:871:U:H4'	36:YQ:69:PHE:CE2	2.53	0.44
25:YA:1048:A:C5	25:YA:1111:A:H2	2.36	0.44
25:YA:1508:A:O2'	25:YA:1509:C:O4'	2.21	0.44
25:YA:1803:A:O3'	27:YD:259:THR:HG21	2.18	0.44
25:YA:2626:C:H2'	25:YA:2627:G:C8	2.53	0.44
25:YA:2747:G:P	31:YH:138:LYS:NZ	2.91	0.44
26:YB:3:C:H2'	26:YB:4:C:C6	2.52	0.44
27:YD:145:VAL:HB	27:YD:155:LEU:HB2	1.99	0.44
27:YD:155:LEU:N	27:YD:155:LEU:HD12	2.32	0.44
28:YE:120:TRP:CE3	28:YE:155:LYS:HD3	2.53	0.44
30:YG:145:THR:O	30:YG:147:ASP:N	2.44	0.44
32:YI:21:VAL:HG22	32:YI:22:LYS:H	1.81	0.44
35:YP:46:LYS:HB3	35:YP:46:LYS:HE3	1.63	0.44
46:Y0:43:THR:O	46:Y0:43:THR:HG23	2.17	0.44
1:QA:176:C:O2'	1:QA:1451:A:N6	2.51	0.44
1:QA:191:G:C1'	20:QT:105:SER:HB3	2.44	0.44
1:QA:455:C:H42	1:QA:477:G:H1	1.66	0.44
1:QA:842:C:H5'	1:QA:843:U:OP1	2.17	0.44
1:QA:1032(A):G:H2'	1:QA:1032(B):G:H8	1.82	0.44
1:QA:1135:U:H4'	1:QA:1136:U:C5	2.39	0.44
1:QA:1333:A:C2	1:QA:1334:G:H1'	2.53	0.44
1:QA:1336:C:O2'	1:QA:1337:G:OP2	2.34	0.44
5:QE:101:ILE:O	5:QE:101:ILE:HG12	2.18	0.44
25:RA:30:G:H2'	25:RA:31:C:C6	2.53	0.44
25:RA:1174:A:H2'	25:RA:1174:A:N3	2.33	0.44
25:RA:1434:A:H61	25:RA:1558:A:H62	1.65	0.44
25:RA:1659:U:O2'	25:RA:2712(A):A:N1	2.45	0.44
25:RA:1956:U:H1'	25:RA:2552:U:OP1	2.16	0.44
25:RA:2419:U:O4	54:R8:30:ARG:CD	2.65	0.44
25:RA:2478:A:H2'	25:RA:2479:G:O4'	2.18	0.44
25:RA:2629:A:H4'	25:RA:2629:A:OP1	2.18	0.44
27:RD:65:ILE:HD13	27:RD:65:ILE:H	1.82	0.44
28:RE:11:MET:O	28:RE:12:THR:HB	2.18	0.44
39:RT:107:ASP:O	39:RT:111:ARG:NH1	2.51	0.44
41:RV:16:PRO:HB3	41:RV:97:LYS:O	2.17	0.44
45:RZ:157:LEU:HD23	45:RZ:161:VAL:HG12	1.98	0.44
46:R0:50:ASN:C	46:R0:62:LEU:HD12	2.42	0.44
51:R5:52:TYR:CD1	51:R5:52:TYR:N	2.85	0.44
54:R8:53:PRO:CD	54:R8:54:GLU:N	2.77	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:96:G:H2'	1:XA:97:U:O4'	2.18	0.44
1:XA:828:A:H2'	1:XA:829:G:O4'	2.17	0.44
1:XA:1003:G:H21	1:XA:1005:A:H5'	1.82	0.44
1:XA:1455:G:H5"	20:XT:31:SER:HB2	2.00	0.44
2:XB:195:ASP:O	8:XH:74:PRO:HG3	2.17	0.44
3:XC:82:GLU:O	3:XC:86:VAL:HG13	2.17	0.44
13:XM:120:LYS:O	13:XM:121:LYS:CB	2.66	0.44
15:XO:26:GLU:H	15:XO:26:GLU:HG2	1.54	0.44
25:YA:270(E):G:H2'	25:YA:270(F):U:O4'	2.17	0.44
25:YA:674:G:O2'	29:YF:74:ARG:HD3	2.18	0.44
25:YA:861:A:C2	25:YA:917:A:C5	3.06	0.44
25:YA:1048:A:OP2	25:YA:1110:G:N2	2.51	0.44
25:YA:1076:C:H2'	25:YA:1077:A:H5"	2.00	0.44
25:YA:2151:G:H2'	25:YA:2152:G:H8	1.82	0.44
27:YD:30:GLU:HG3	27:YD:63:ARG:NE	2.32	0.44
29:YF:184:TYR:CE2	29:YF:188:ARG:HD2	2.52	0.44
30:YG:98:ARG:O	30:YG:101:ILE:HG13	2.17	0.44
31:YH:6:ARG:CG	31:YH:7:LEU:N	2.81	0.44
31:YH:37:VAL:HG11	31:YH:68:THR:HG23	1.98	0.44
32:YI:20:ASP:OD1	32:YI:20:ASP:N	2.45	0.44
32:YI:52:ARG:O	32:YI:56:LYS:HB3	2.18	0.44
36:YQ:60:ARG:HB2	36:YQ:60:ARG:HH21	1.83	0.44
36:YQ:80:GLU:OE1	46:Y0:6:GLY:O	2.36	0.44
38:YS:112:PHE:O	38:YS:112:PHE:CD1	2.70	0.44
39:YT:48:ILE:HD12	39:YT:48:ILE:H	1.83	0.44
41:YV:15:GLU:HG3	41:YV:16:PRO:HD2	1.99	0.44
44:YY:51:VAL:O	44:YY:56:PRO:HA	2.18	0.44
50:Y4:48:ARG:CZ	50:Y4:51:ASP:HA	2.47	0.44
1:QA:446:G:H2'	1:QA:447:G:O4'	2.18	0.44
1:QA:1316:G:H5"	14:QN:17:LYS:HE3	2.00	0.44
4:QD:63:LYS:O	4:QD:67:ILE:HG13	2.17	0.44
4:QD:63:LYS:HD2	4:QD:198:VAL:HG22	2.00	0.44
4:QD:129:ASN:HA	4:QD:145:GLU:HB2	2.00	0.44
16:QP:23:ASP:O	16:QP:26:ARG:HB2	2.17	0.44
25:RA:48:G:N2	25:RA:177:G:H21	2.16	0.44
25:RA:270(I):G:H2'	25:RA:270(J):G:C8	2.53	0.44
25:RA:1319:G:C6	25:RA:1320:C:N4	2.86	0.44
25:RA:1729:A:H2'	25:RA:1730:U:H6	1.83	0.44
25:RA:2392:A:O2'	54:R8:27:THR:HB	2.17	0.44
25:RA:2470:G:H5'	36:RQ:56:ARG:HH22	1.83	0.44
27:RD:34:VAL:HG22	27:RD:35:LYS:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:RG:57:ALA:HB1	30:RG:68:PRO:HG2	1.99	0.44
30:RG:171:ALA:O	30:RG:175:LEU:HG	2.18	0.44
31:RH:58:GLU:O	31:RH:59:ARG:C	2.61	0.44
36:RQ:20:ALA:C	36:RQ:21:THR:HG23	2.43	0.44
36:RQ:34:LEU:HB2	36:RQ:118:LEU:HD22	1.99	0.44
36:RQ:133:ARG:CG	36:RQ:134:ARG:N	2.78	0.44
42:RW:1:MET:HE3	42:RW:2:GLU:H	1.82	0.44
43:RX:87:GLN:O	43:RX:88:LYS:HG3	2.18	0.44
44:RY:81:LYS:HB2	44:RY:96:ILE:CG2	2.48	0.44
45:RZ:151:HIS:HA	45:RZ:170:THR:HA	1.99	0.44
50:R4:50:VAL:O	50:R4:50:VAL:CG1	2.63	0.44
52:R6:40:CYS:HA	52:R6:41:PRO:HD2	1.85	0.44
54:R8:58:ILE:O	54:R8:61:LEU:CG	2.66	0.44
1:XA:253:U:H2'	1:XA:254:G:C8	2.53	0.44
1:XA:953:G:C5'	1:XA:965:A:H61	2.29	0.44
3:XC:47:LEU:HD12	3:XC:47:LEU:HA	1.82	0.44
5:XE:82:VAL:HG21	5:XE:138:ALA:HA	2.00	0.44
10:XJ:58:ASP:O	10:XJ:59:SER:HB2	2.17	0.44
11:XK:19:ALA:HB2	11:XK:32:ILE:HG22	2.00	0.44
13:XM:23:TYR:HE2	13:XM:70:LEU:HD12	1.83	0.44
15:XO:32:LEU:O	15:XO:36:ILE:HG13	2.18	0.44
15:XO:39:LEU:HD13	15:XO:56:LEU:HB2	2.00	0.44
15:XO:77:ARG:HA	15:XO:80:ALA:HB3	1.99	0.44
16:XP:4:ILE:HB	16:XP:66:PRO:HB3	2.00	0.44
16:XP:39:TYR:CZ	16:XP:41:PRO:HB3	2.53	0.44
25:YA:184:C:H2'	25:YA:185:U:C6	2.53	0.44
25:YA:814:C:H41	35:YP:25:SER:HA	1.83	0.44
25:YA:953:A:OP1	36:YQ:18:LYS:NZ	2.49	0.44
25:YA:1300:U:H4'	25:YA:1301:A:H5'	2.00	0.44
25:YA:1788:C:OP1	27:YD:222:ARG:NH2	2.42	0.44
26:YB:24:G:H5"	26:YB:25:A:OP1	2.17	0.44
27:YD:44:ASN:HB2	27:YD:49:ILE:HA	1.93	0.44
27:YD:102:LYS:O	27:YD:103:ARG:CG	2.66	0.44
28:YE:2:LYS:HG2	28:YE:95:ILE:HG22	1.99	0.44
29:YF:65:TRP:CZ2	29:YF:72:ARG:NH2	2.86	0.44
30:YG:10:LYS:HE2	30:YG:175:LEU:O	2.18	0.44
38:YS:14:VAL:CG1	38:YS:15:ARG:N	2.81	0.44
47:Y1:70:VAL:O	47:Y1:73:LEU:HB2	2.18	0.44
49:Y3:31:LEU:O	49:Y3:32:GLN:HB2	2.17	0.44
52:Y6:41:PRO:HD2	52:Y6:46:HIS:H	1.81	0.44
1:QA:429:U:H3'	4:QD:22:LYS:NZ	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:751:U:H2'	1:QA:752:G:O4'	2.18	0.44
1:QA:947:G:O3'	13:QM:109:THR:OG1	2.35	0.44
9:QI:16:ARG:O	9:QI:63:ILE:HA	2.17	0.44
10:QJ:47:PHE:HE1	10:QJ:63:PHE:HB2	1.83	0.44
22:QV:53:G:O6	22:QV:62:C:N4	2.50	0.44
25:RA:1165:U:H2'	25:RA:1166:C:C6	2.53	0.44
25:RA:1788:C:H2'	25:RA:1789:A:O4'	2.17	0.44
25:RA:1798:U:C5'	27:RD:259:THR:HG22	2.47	0.44
25:RA:1930:G:O2'	25:RA:1931:U:OP2	2.36	0.44
26:RB:104:A:H2'	26:RB:105:G:O4'	2.18	0.44
28:RE:13:ARG:HH11	28:RE:13:ARG:HB3	1.82	0.44
28:RE:199:ARG:HH11	28:RE:199:ARG:HG3	1.82	0.44
38:RS:19:LYS:O	38:RS:20:ARG:HB3	2.18	0.44
43:RX:67:GLY:O	43:RX:69:TYR:N	2.43	0.44
52:R6:41:PRO:HD2	52:R6:46:HIS:H	1.83	0.44
3:XC:72:LYS:HB3	3:XC:75:VAL:HG23	2.00	0.44
3:XC:108:ASN:HB3	3:XC:111:LEU:HD12	2.00	0.44
4:XD:33:MET:HE2	4:XD:37:PRO:CA	2.44	0.44
5:XE:9:LYS:HB3	5:XE:112:LEU:HD11	2.00	0.44
12:XL:120:TYR:O	12:XL:121:GLY:O	2.36	0.44
25:YA:1087:G:C4	25:YA:1089:G:H1'	2.53	0.44
25:YA:1165:U:H2'	25:YA:1166:C:C6	2.53	0.44
25:YA:1173:G:H5"	25:YA:1174:A:OP1	2.18	0.44
25:YA:1476:C:H2'	25:YA:1477:A:H8	1.83	0.44
27:YD:17:THR:HG21	27:YD:204:ILE:HA	1.99	0.44
27:YD:44:ASN:CB	27:YD:49:ILE:HG22	2.46	0.44
29:YF:108:LYS:O	29:YF:109:GLY:C	2.61	0.44
30:YG:83:ARG:HG3	30:YG:86:MET:HE3	1.99	0.44
31:YH:153:LYS:HG3	31:YH:162:ILE:H	1.78	0.44
32:YI:30:LEU:HB3	32:YI:36:ALA:HB3	2.00	0.44
36:YQ:81:VAL:HG23	36:YQ:82:ARG:N	2.32	0.44
38:YS:83:LYS:O	38:YS:109:GLY:CA	2.46	0.44
45:YZ:100:VAL:HA	45:YZ:101:PRO:HD3	1.86	0.44
1:QA:255:G:H2'	1:QA:256:U:C6	2.53	0.43
1:QA:261:U:O2	1:QA:263:A:C8	2.71	0.43
1:QA:452:A:O2'	1:QA:453:A:O5'	2.36	0.43
1:QA:743:U:H2'	1:QA:744:C:C6	2.53	0.43
1:QA:1179:A:O3'	9:QI:103:THR:HG23	2.18	0.43
1:QA:1255:G:C6	1:QA:1279:A:N7	2.86	0.43
1:QA:1271:G:C6	1:QA:1272:G:C6	3.05	0.43
2:QB:37:ASN:C	2:QB:39:ILE:H	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:QB:74:LYS:O	2:QB:78:GLN:HG3	2.18	0.43
3:QC:81:GLY:O	3:QC:85:ARG:HB2	2.18	0.43
7:QG:113:GLU:H	7:QG:113:GLU:HG2	1.39	0.43
10:QJ:51:ARG:NE	10:QJ:60:ARG:O	2.45	0.43
25:RA:242:G:C3'	54:R8:6:THR:HG23	2.47	0.43
25:RA:746:A:C5	25:RA:2611:U:H5"	2.53	0.43
25:RA:1727:U:H2'	25:RA:1728:G:O4'	2.18	0.43
25:RA:1790:C:H2'	25:RA:1791:A:C5	2.52	0.43
25:RA:2790:A:H2'	25:RA:2791:C:H5"	2.00	0.43
26:RB:50:G:H5"	38:RS:61:ASN:ND2	2.32	0.43
27:RD:145:VAL:HG11	27:RD:175:LEU:HD11	2.00	0.43
28:RE:52:LEU:HB2	28:RE:75:VAL:CG2	2.40	0.43
31:RH:125:VAL:CG1	31:RH:126:PRO:CG	2.94	0.43
34:RO:22:ILE:HD13	34:RO:22:ILE:HA	1.77	0.43
37:RR:27:SER:HB3	37:RR:34:ILE:HD11	2.00	0.43
37:RR:54:LEU:HD23	37:RR:66:VAL:HG23	1.99	0.43
54:R8:40:GLU:O	54:R8:43:GLN:N	2.50	0.43
1:XA:652:U:HO2'	1:XA:653:A:P	2.41	0.43
1:XA:1316:G:N2	1:XA:1318:A:H3'	2.33	0.43
1:XA:1434:A:H2'	1:XA:1435:G:O4'	2.18	0.43
4:XD:100:ARG:NH1	4:XD:137:SER:HB3	2.33	0.43
12:XL:6:THR:H	12:XL:9:GLN:NE2	1.97	0.43
16:XP:45:THR:HG22	16:XP:47:ASP:N	2.26	0.43
25:YA:710:G:H2'	25:YA:711:G:C8	2.53	0.43
25:YA:1049:C:H2'	25:YA:1050:A:H5"	1.99	0.43
25:YA:1265:A:H3'	51:Y5:19:ARG:NH1	2.33	0.43
25:YA:1316:U:H2'	25:YA:1317:A:C8	2.52	0.43
25:YA:2477:C:H1'	25:YA:2481:G:O6	2.18	0.43
25:YA:2477:C:H2'	55:Y9:1:MET:HG3	2.00	0.43
27:YD:95:LEU:HD12	27:YD:95:LEU:O	2.17	0.43
27:YD:143:HIS:HD2	27:YD:144:ALA:HB2	1.82	0.43
28:YE:48:GLN:HE21	28:YE:48:GLN:HB3	1.55	0.43
29:YF:42:ALA:O	29:YF:45:ARG:HB2	2.18	0.43
29:YF:174:VAL:CG1	29:YF:174:VAL:O	2.65	0.43
30:YG:67:LYS:O	30:YG:67:LYS:HD2	2.17	0.43
38:YS:57:LYS:O	38:YS:58:LEU:HB3	2.18	0.43
43:YX:70:LEU:HD23	43:YX:70:LEU:H	1.83	0.43
1:QA:646:U:H2'	1:QA:647:C:C6	2.53	0.43
1:QA:1398:A:H5"	1:QA:1401:G:H4'	1.98	0.43
4:QD:33:MET:HE2	4:QD:37:PRO:HA	2.00	0.43
5:QE:69:VAL:HG11	5:QE:71:LEU:HD21	1.98	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:QG:9:VAL:HG13	7:QG:94:ARG:NH2	2.27	0.43
7:QG:116:ALA:HA	7:QG:119:ARG:HE	1.83	0.43
12:QL:27:LEU:HD13	12:QL:28:LYS:H	1.83	0.43
12:QL:119:LYS:HB2	12:QL:120:TYR:HD1	1.83	0.43
22:QV:3:G:O2'	22:QV:4:G:H8	2.01	0.43
25:RA:29:U:H2'	25:RA:30:G:C8	2.53	0.43
25:RA:31:C:O3'	25:RA:1238:G:H5'	2.18	0.43
25:RA:208:C:H2'	25:RA:209:C:C6	2.53	0.43
25:RA:270(I):G:H1	25:RA:270(Q):C:H42	1.67	0.43
25:RA:796:C:H2'	25:RA:797:C:C6	2.53	0.43
25:RA:1316:U:H2'	25:RA:1317:A:C8	2.53	0.43
27:RD:70:TRP:CD1	27:RD:70:TRP:C	2.96	0.43
28:RE:3:GLY:CA	28:RE:81:ILE:HG21	2.48	0.43
28:RE:15:PHE:CE1	39:RT:81:PRO:CD	3.01	0.43
28:RE:36:ARG:HB3	28:RE:36:ARG:NH1	2.30	0.43
28:RE:51:PHE:CD2	28:RE:52:LEU:N	2.76	0.43
30:RG:6:ALA:CA	50:R4:23:GLU:HG2	2.47	0.43
31:RH:137:ASP:OD1	31:RH:138:LYS:N	2.51	0.43
32:RI:95:LYS:HA	32:RI:111:PRO:HG3	2.00	0.43
34:RO:20:MET:HE3	34:RO:44:LYS:HE3	2.01	0.43
44:RY:42:VAL:O	44:RY:65:ALA:N	2.45	0.43
45:RZ:153:SER:HB2	45:RZ:167:PRO:HB3	2.00	0.43
52:R6:28:ARG:HG3	52:R6:31:PRO:HD2	2.00	0.43
1:XA:730:G:C5	1:XA:731:G:H1'	2.53	0.43
1:XA:1108:G:H5'	3:XC:176:HIS:ND1	2.27	0.43
4:XD:153:ARG:NH1	4:XD:181:MET:HB2	2.32	0.43
6:XF:33:TYR:HB2	6:XF:75:LEU:HD12	1.99	0.43
9:XI:118:LYS:O	9:XI:119:ALA:HB3	2.18	0.43
10:XJ:57:LYS:CD	10:XJ:60:ARG:NH2	2.78	0.43
25:YA:239:U:H2'	25:YA:240:G:O4'	2.17	0.43
25:YA:530:G:C5	25:YA:2022:U:H5"	2.53	0.43
25:YA:709:U:H2'	25:YA:710:G:C8	2.53	0.43
25:YA:871:U:H4'	36:YQ:69:PHE:CD2	2.53	0.43
25:YA:1045:A:C8	25:YA:1111:A:N6	2.86	0.43
25:YA:1688:U:H1'	25:YA:1701:A:C6	2.53	0.43
27:YD:35:LYS:HB3	27:YD:36:PRO:HA	2.00	0.43
29:YF:79:GLY:O	29:YF:80:ALA:C	2.61	0.43
31:YH:35:VAL:CG2	31:YH:75:ALA:HB2	2.48	0.43
31:YH:137:ASP:OD1	31:YH:138:LYS:N	2.51	0.43
32:YI:144:VAL:HG22	32:YI:145:VAL:H	1.83	0.43
33:YN:112:LEU:O	33:YN:112:LEU:HG	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:YS:38:GLN:CG	38:YS:47:THR:HG21	2.48	0.43
42:YW:14:PRO:C	42:YW:18:ARG:HD2	2.43	0.43
51:Y5:56:LYS:CG	51:Y5:58:LEU:HB3	2.46	0.43
54:Y8:58:ILE:O	54:Y8:61:LEU:CG	2.66	0.43
55:Y9:1:MET:O	55:Y9:34:GLN:HG2	2.18	0.43
1:QA:109:A:H5'	1:QA:110:C:C5	2.54	0.43
1:QA:636:U:H2'	1:QA:637:G:H8	1.83	0.43
1:QA:967:C:H2'	1:QA:968:A:N7	2.33	0.43
1:QA:984:C:H2'	1:QA:985:C:H6	1.83	0.43
1:QA:1347:G:O2'	1:QA:1348:U:OP2	2.29	0.43
2:QB:167:PRO:HG3	2:QB:188:ALA:HB2	2.00	0.43
11:QK:19:ALA:HB2	11:QK:32:ILE:HG22	2.00	0.43
15:QO:17:ARG:HD3	15:QO:26:GLU:HG3	1.99	0.43
17:QQ:60:ILE:HB	17:QQ:74:LEU:HD23	2.00	0.43
25:RA:242:G:H5'	54:R8:62:LEU:CD2	2.43	0.43
25:RA:520:G:H2'	25:RA:521:G:C8	2.53	0.43
25:RA:2415:G:H4'	35:RP:67:MET:N	2.33	0.43
25:RA:2517:C:C2	25:RA:2542:A:N6	2.86	0.43
27:RD:30:GLU:CD	27:RD:63:ARG:HH21	2.26	0.43
28:RE:16:ARG:O	28:RE:18:ASP:O	2.36	0.43
31:RH:92:ILE:CD1	31:RH:160:LYS:HD3	2.48	0.43
31:RH:136:ILE:HD12	31:RH:136:ILE:N	2.31	0.43
32:RI:9:LEU:O	32:RI:10:GLU:HG3	2.18	0.43
32:RI:52:ARG:HA	32:RI:56:LYS:HB3	1.99	0.43
33:RN:33:LEU:HA	33:RN:38:HIS:CE1	2.53	0.43
35:RP:62:LEU:CD2	54:R8:25:MET:CB	2.78	0.43
35:RP:65:ARG:NH2	54:R8:46:ARG:NH1	2.66	0.43
37:RR:98:LEU:HB3	51:R5:45:VAL:HG23	2.00	0.43
44:RY:47:LYS:O	44:RY:49:VAL:N	2.51	0.43
44:RY:97:ARG:HE	44:RY:98:VAL:HB	1.83	0.43
50:R4:15:ILE:N	50:R4:15:ILE:CD1	2.78	0.43
50:R4:35:VAL:C	50:R4:37:SER:H	2.19	0.43
50:R4:39:CYS:O	50:R4:40:HIS:CB	2.66	0.43
54:R8:58:ILE:O	54:R8:61:LEU:CD1	2.67	0.43
1:XA:300:A:H8	1:XA:300:A:O5'	2.01	0.43
1:XA:539:A:H2'	1:XA:540:G:C8	2.53	0.43
1:XA:1216:G:H5"	14:XN:5:ALA:HB2	2.00	0.43
1:XA:1280:A:H1'	10:XJ:41:PRO:HG3	2.01	0.43
1:XA:1435:G:H1	1:XA:1466:C:H42	1.65	0.43
2:XB:80:ILE:HG21	2:XB:212:GLN:HA	2.00	0.43
10:XJ:32:ALA:H	10:XJ:78:ASN:ND2	2.16	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:XM:14:ARG:H	13:XM:44:ARG:CD	2.25	0.43
25:YA:467:G:OP2	53:Y7:34:ARG:NH1	2.50	0.43
25:YA:605:C:H1'	25:YA:657:U:O2'	2.18	0.43
25:YA:1021:A:OP2	33:YN:65:LYS:NZ	2.51	0.43
25:YA:1530:G:O6	25:YA:1542:G:N2	2.51	0.43
25:YA:1534:G:O2'	25:YA:1535:U:H4'	2.19	0.43
25:YA:1541:U:H2'	25:YA:1542:G:O4'	2.18	0.43
25:YA:1916:A:H2'	25:YA:1917:U:O4'	2.19	0.43
25:YA:2306:C:H3'	25:YA:2307:G:C5'	2.44	0.43
25:YA:2532:G:H1'	25:YA:2663:G:N2	2.33	0.43
25:YA:2688:U:H1'	25:YA:2721:A:H61	1.84	0.43
27:YD:145:VAL:HG12	27:YD:146:GLU:N	2.32	0.43
28:YE:3:GLY:HA3	28:YE:81:ILE:CD1	2.48	0.43
28:YE:3:GLY:CA	28:YE:81:ILE:HG21	2.49	0.43
32:YI:48:GLU:OE1	32:YI:52:ARG:NH2	2.51	0.43
36:YQ:21:THR:HB	36:YQ:22:LYS:H	1.42	0.43
36:YQ:80:GLU:C	36:YQ:81:VAL:CG1	2.88	0.43
37:YR:38:VAL:HG22	37:YR:112:ALA:HB2	2.00	0.43
41:YV:52:VAL:O	41:YV:54:GLY:N	2.51	0.43
43:YX:72:LYS:HG2	43:YX:73:ARG:O	2.18	0.43
47:Y1:83:GLU:C	47:Y1:85:LEU:N	2.76	0.43
47:Y1:94:LEU:HD23	47:Y1:94:LEU:HA	1.82	0.43
54:Y8:58:ILE:O	54:Y8:61:LEU:CD1	2.67	0.43
1:QA:1028:C:N4	1:QA:1033:G:H1	2.16	0.43
4:QD:21:LEU:HD21	4:QD:67:ILE:HA	2.00	0.43
12:QL:44:THR:HA	12:QL:45:PRO:HD3	1.70	0.43
16:QP:20:VAL:HG21	16:QP:32:TYR:CE2	2.54	0.43
18:QR:37:VAL:HG22	18:QR:78:LEU:HB3	2.00	0.43
25:RA:270(Y):G:H4'	25:RA:273:G:O3'	2.18	0.43
25:RA:527:C:N3	25:RA:2779:U:H5"	2.33	0.43
25:RA:706:A:H2'	25:RA:707:G:O4'	2.18	0.43
25:RA:892:G:N2	25:RA:893:C:O2	2.51	0.43
25:RA:1657:C:H2'	25:RA:1658:C:H6	1.83	0.43
25:RA:2415:G:H4'	35:RP:66:GLY:HA3	2.00	0.43
25:RA:2611:U:OP2	25:RA:2611:U:H6	2.02	0.43
25:RA:2817:G:C4	25:RA:2830:G:N2	2.87	0.43
28:RE:16:ARG:C	28:RE:17:ASP:CG	2.86	0.43
28:RE:31:CYS:HB3	28:RE:49:LEU:HG	2.01	0.43
28:RE:52:LEU:O	28:RE:74:PRO:HA	2.18	0.43
29:RF:150:GLY:HA2	29:RF:172:TRP:CD2	2.52	0.43
36:RQ:27:VAL:HG13	36:RQ:28:ALA:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:RT:19:LEU:HA	39:RT:20:PRO:HD3	1.86	0.43
43:RX:40:LYS:C	43:RX:42:ALA:N	2.76	0.43
44:RY:46:LYS:HB2	44:RY:61:ILE:HG22	1.99	0.43
1:XA:247:G:O6	1:XA:278:G:C6	2.71	0.43
1:XA:427:U:OP1	4:XD:13:ARG:NH2	2.52	0.43
1:XA:554:C:H2'	1:XA:555:C:H6	1.83	0.43
1:XA:633:G:H5'	1:XA:634:C:OP2	2.18	0.43
12:XL:27:LEU:HD13	12:XL:28:LYS:H	1.84	0.43
25:YA:464:U:H2'	25:YA:465:G:O4'	2.17	0.43
25:YA:574:C:O2	28:YE:145:LYS:HE3	2.19	0.43
25:YA:896:A:N3	45:YZ:176:PRO:HB3	2.33	0.43
25:YA:1797:C:O2'	27:YD:259:THR:HB	2.18	0.43
25:YA:2360:A:OP1	54:Y8:49:VAL:HA	2.18	0.43
25:YA:2360:A:OP1	54:Y8:49:VAL:HG12	2.18	0.43
25:YA:2364:C:H2'	25:YA:2365:G:O4'	2.19	0.43
25:YA:2887:U:H2'	25:YA:2888:C:C6	2.53	0.43
27:YD:43:ARG:CZ	27:YD:49:ILE:HG21	2.49	0.43
27:YD:181:GLU:HA	27:YD:272:ALA:CB	2.39	0.43
28:YE:11:MET:O	28:YE:12:THR:HB	2.18	0.43
28:YE:143:ASN:N	28:YE:143:ASN:ND2	2.65	0.43
31:YH:136:ILE:HD12	31:YH:136:ILE:N	2.31	0.43
32:YI:123:LEU:HD23	32:YI:142:VAL:HB	1.99	0.43
35:YP:115:LEU:HA	35:YP:134:ALA:HB2	2.00	0.43
38:YS:99:LYS:C	38:YS:101:LEU:H	2.25	0.43
39:YT:35:LYS:H	39:YT:35:LYS:HD2	1.83	0.43
48:Y2:59:ARG:O	48:Y2:62:THR:HG23	2.18	0.43
54:Y8:40:GLU:O	54:Y8:43:GLN:N	2.50	0.43
1:QA:28:G:O2'	1:QA:296:U:OP1	2.34	0.43
1:QA:654:G:C2	1:QA:753:A:C4	3.06	0.43
1:QA:953:G:H5'	1:QA:965:A:H61	1.82	0.43
1:QA:978:A:O2'	1:QA:1322:C:N3	2.42	0.43
2:QB:8:LYS:HE3	2:QB:11:LEU:HB3	2.01	0.43
8:QH:13:ILE:O	8:QH:17:THR:HG23	2.19	0.43
11:QK:38:ASN:HA	11:QK:39:PRO:HD3	1.88	0.43
13:QM:122:LYS:HD3	13:QM:122:LYS:C	2.42	0.43
22:QV:23:C:H2'	22:QV:24:U:C6	2.52	0.43
25:RA:107:C:H2'	25:RA:108:U:H6	1.84	0.43
25:RA:265:A:H2'	25:RA:266:G:O4'	2.18	0.43
25:RA:273(C):C:N4	25:RA:363(C):G:H1	2.14	0.43
25:RA:859:G:H2'	25:RA:916:G:O6	2.17	0.43
25:RA:1309:G:P	53:R7:9:ARG:HD3	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1687:G:N1	25:RA:1700:A:OP1	2.49	0.43
25:RA:2197:U:H1'	25:RA:2198:A:C8	2.53	0.43
25:RA:2662:A:H8	25:RA:2662:A:O5'	2.01	0.43
25:RA:2867:G:HO2'	25:RA:2868:A:P	2.42	0.43
26:RB:3:C:H2'	26:RB:4:C:C6	2.54	0.43
26:RB:14:U:H5'	26:RB:71:C:O4'	2.19	0.43
28:RE:48:GLN:HE21	28:RE:48:GLN:HB3	1.55	0.43
50:R4:49:PHE:N	50:R4:49:PHE:HD1	2.17	0.43
1:XA:537:G:H5"	12:XL:113:ARG:HH12	1.83	0.43
1:XA:939:G:H5"	7:XG:102:ARG:NH2	2.33	0.43
1:XA:1054:C:C2	24:XY:34:G:O4'	2.71	0.43
2:XB:100:GLY:N	2:XB:176:GLU:OE2	2.47	0.43
5:XE:102:ALA:CB	5:XE:120:THR:OG1	2.65	0.43
10:XJ:44:VAL:HG13	10:XJ:66:ARG:HG2	1.99	0.43
17:XQ:62:SER:HB3	17:XQ:72:ARG:HE	1.84	0.43
25:YA:102:G:H4'	25:YA:103:A:O5'	2.18	0.43
25:YA:508:G:O2'	25:YA:509:C:P	2.77	0.43
25:YA:922:U:H2'	25:YA:923:C:C6	2.54	0.43
25:YA:1803:A:O3'	27:YD:259:THR:CG2	2.67	0.43
25:YA:2105:C:H2'	25:YA:2106:G:C8	2.53	0.43
25:YA:2682:U:O2'	28:YE:13:ARG:HG2	2.19	0.43
28:YE:51:PHE:O	28:YE:74:PRO:CB	2.67	0.43
30:YG:47:LYS:HB2	30:YG:47:LYS:HE3	1.73	0.43
31:YH:27:LYS:HD2	31:YH:27:LYS:C	2.44	0.43
31:YH:59:ARG:CG	31:YH:59:ARG:NH1	2.79	0.43
31:YH:149:ARG:HA	31:YH:162:ILE:HG21	1.99	0.43
32:YI:2:LYS:HA	32:YI:20:ASP:HA	2.00	0.43
32:YI:93:THR:HG22	32:YI:119:PRO:HB3	2.01	0.43
34:YO:64:ARG:HG2	34:YO:79:PHE:CD1	2.53	0.43
34:YO:88:ASN:OD1	34:YO:90:GLN:HB2	2.19	0.43
37:YR:34:ILE:HA	37:YR:34:ILE:HD13	1.71	0.43
38:YS:86:ALA:O	38:YS:87:PHE:CB	2.65	0.43
41:YV:81:TYR:C	41:YV:82:ARG:HD2	2.43	0.43
43:YX:53:LYS:H	43:YX:82:GLN:HB3	1.83	0.43
44:YY:67:LEU:HD12	44:YY:67:LEU:HA	1.78	0.43
1:QA:189:U:HO2'	1:QA:190:G:P	2.41	0.43
1:QA:266:G:H5"	1:QA:267:C:H5	1.83	0.43
1:QA:686:U:O4	1:QA:703:G:H1'	2.18	0.43
1:QA:867:G:O2'	1:QA:873:A:N1	2.45	0.43
1:QA:1336:C:O2'	1:QA:1337:G:P	2.76	0.43
1:QA:1432:G:OP1	39:RT:107:ASP:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:QD:157:LEU:O	4:QD:161:ASN:ND2	2.43	0.43
18:QR:53:ARG:HH21	18:QR:60:ALA:N	2.17	0.43
20:QT:84:LEU:HD23	20:QT:84:LEU:HA	1.86	0.43
25:RA:146:G:H2'	25:RA:147:U:O4'	2.18	0.43
25:RA:590:A:H2'	25:RA:591:C:C6	2.53	0.43
25:RA:1657:C:OP1	28:RE:136:ARG:N	2.52	0.43
25:RA:1849:G:H2'	25:RA:1850:G:C8	2.54	0.43
28:RE:51:PHE:O	28:RE:74:PRO:CB	2.67	0.43
28:RE:120:TRP:CE3	28:RE:155:LYS:HD3	2.53	0.43
32:RI:129:THR:HA	32:RI:137:PRO:HA	1.99	0.43
39:RT:1:MET:O	39:RT:3:ARG:HG2	2.19	0.43
50:R4:59:PHE:CE1	50:R4:70:GLY:N	2.86	0.43
1:XA:35:G:N2	12:XL:118:SER:OG	2.48	0.43
1:XA:690:G:N1	1:XA:691:G:C6	2.86	0.43
2:XB:215:LEU:HD22	2:XB:215:LEU:HA	1.73	0.43
12:XL:76:ASN:O	12:XL:76:ASN:CG	2.62	0.43
13:XM:121:LYS:HZ1	24:XY:40:G:H4'	1.84	0.43
14:XN:27:CYS:SG	14:XN:29:ARG:HB2	2.58	0.43
14:XN:29:ARG:HD3	14:XN:40:CYS:HB2	1.99	0.43
15:XO:25:THR:HG21	15:XO:70:LEU:HB2	2.01	0.43
22:XV:2:C:H2'	22:XV:3:G:H5'	1.99	0.43
22:XV:19:G:H5'	22:XV:20:U:C5	2.52	0.43
25:YA:142:G:H1'	43:YX:37:THR:HG21	2.00	0.43
25:YA:587:C:N3	35:YP:33:ARG:NH1	2.66	0.43
26:YB:38:C:N4	26:YB:44:G:H1	2.16	0.43
27:YD:11:PRO:O	27:YD:12:SER:OG	2.29	0.43
28:YE:155:LYS:O	28:YE:156:MET:HG3	2.19	0.43
31:YH:92:ILE:CD1	31:YH:160:LYS:HD3	2.48	0.43
34:YO:88:ASN:ND2	34:YO:92:GLU:HB2	2.22	0.43
38:YS:88:ASP:CG	38:YS:89:ARG:N	2.77	0.43
40:YU:60:LEU:HD11	40:YU:64:ARG:CZ	2.48	0.43
45:YZ:30:ASN:OD1	45:YZ:33:LEU:N	2.49	0.43
48:Y2:41:ILE:C	48:Y2:41:ILE:CD1	2.81	0.43
49:Y3:51:ALA:HA	49:Y3:54:VAL:HG12	2.00	0.43
53:Y7:47:ARG:HB2	53:Y7:48:LYS:H	1.60	0.43
54:Y8:28:GLY:O	54:Y8:29:LYS:C	2.61	0.43
1:QA:464:G:O6	1:QA:466:C:H5'	2.18	0.43
1:QA:520:A:N1	1:QA:536:C:H1'	2.33	0.43
1:QA:640:A:O2'	8:QH:115:SER:HB2	2.17	0.43
1:QA:1124:G:H5"	1:QA:1145:C:H41	1.83	0.43
1:QA:1157:A:H1'	1:QA:1158:C:C4	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:QE:101:ILE:HG13	5:QE:119:LEU:HD23	2.01	0.43
13:QM:44:ARG:HB2	13:QM:47:ASP:OD2	2.19	0.43
20:QT:16:HIS:O	20:QT:19:SER:HB3	2.18	0.43
22:QV:19:G:H4'	22:QV:20:U:OP2	2.19	0.43
25:RA:414:C:H2'	25:RA:415:A:C8	2.53	0.43
25:RA:1005:C:O2'	33:RN:28:THR:HG21	2.18	0.43
31:RH:6:ARG:CG	31:RH:7:LEU:N	2.81	0.43
32:RI:40:THR:O	32:RI:44:LEU:N	2.46	0.43
42:RW:20:VAL:HG22	42:RW:47:VAL:HG21	2.00	0.43
44:RY:55:TYR:CD1	44:RY:55:TYR:N	2.86	0.43
45:RZ:5:LEU:HD21	45:RZ:44:PHE:HA	2.01	0.43
48:R2:39:ALA:HA	48:R2:45:SER:HB2	2.00	0.43
1:XA:657:G:C2	1:XA:658:G:C8	3.06	0.43
1:XA:966:G:O2'	9:XI:127:LYS:O	2.37	0.43
1:XA:1055:A:H2'	3:XC:156:ARG:HD2	2.01	0.43
1:XA:1058:G:H2'	1:XA:1059:C:O4'	2.18	0.43
1:XA:1402:C:H2'	1:XA:1403:C:O4'	2.18	0.43
19:XS:42:PRO:HD3	50:Y4:63:TYR:CE1	2.53	0.43
19:XS:68:GLY:HA3	50:Y4:68:ARG:HB2	1.99	0.43
20:XT:44:ALA:HB1	20:XT:92:LEU:HG	2.00	0.43
25:YA:26:G:C6	25:YA:27:G:N1	2.87	0.43
25:YA:84:A:N1	25:YA:98:G:O2'	2.39	0.43
25:YA:363(B):G:H2'	25:YA:363(C):G:H8	1.84	0.43
25:YA:389:G:H22	35:YP:72:PRO:CG	2.31	0.43
25:YA:732:C:H2'	25:YA:733:G:O4'	2.18	0.43
25:YA:846:C:C2	25:YA:847:U:H5	2.37	0.43
25:YA:1061:U:H4'	25:YA:1070:A:H1'	2.00	0.43
25:YA:1418:G:O5'	25:YA:1418:G:H8	2.02	0.43
25:YA:1509:C:H2'	25:YA:1511:A:C8	2.54	0.43
25:YA:1818:U:C2'	27:YD:157:ARG:HG3	2.48	0.43
25:YA:2360:A:C2	25:YA:2361:A:H1'	2.54	0.43
25:YA:2389:G:H5"	25:YA:2390:U:O4'	2.19	0.43
27:YD:44:ASN:HB3	27:YD:49:ILE:CG2	2.47	0.43
28:YE:16:ARG:O	28:YE:18:ASP:O	2.36	0.43
35:YP:64:LYS:O	35:YP:66:GLY:N	2.52	0.43
36:YQ:27:VAL:HG13	36:YQ:28:ALA:N	2.32	0.43
37:YR:70:LEU:HD23	37:YR:70:LEU:HA	1.84	0.43
38:YS:30:ARG:NH2	38:YS:92:TYR:HD1	2.17	0.43
40:YU:92:ARG:HD2	41:YV:11:GLN:CB	2.47	0.43
41:YV:52:VAL:C	41:YV:54:GLY:H	2.26	0.43
41:YV:55:ALA:HB2	41:YV:101:GLY:HA2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:YY:87:LYS:HB2	44:YY:87:LYS:NZ	2.34	0.43
45:YZ:112:ARG:O	45:YZ:114:GLY:N	2.52	0.43
48:Y2:62:THR:O	48:Y2:65:ASN:HB2	2.18	0.43
1:QA:45:U:OP1	1:QA:307:C:O2'	2.34	0.43
1:QA:376:G:H5"	16:QP:5:ARG:HD2	2.00	0.43
1:QA:936:C:H2'	1:QA:937:A:O4'	2.18	0.43
1:QA:967:C:H2'	1:QA:968:A:C8	2.54	0.43
1:QA:1064:G:O2'	1:QA:1065:U:O5'	2.32	0.43
2:QB:217:ARG:HE	2:QB:217:ARG:HB2	1.29	0.43
5:QE:69:VAL:HG11	5:QE:71:LEU:CD2	2.48	0.43
8:QH:91:ARG:HB2	12:QL:7:ILE:HG13	1.99	0.43
9:QI:112:LYS:HD3	9:QI:113:LYS:O	2.18	0.43
9:QI:116:LYS:HE2	9:QI:122:ALA:HB2	2.01	0.43
18:QR:29:PHE:N	18:QR:29:PHE:CD1	2.87	0.43
25:RA:123:G:O3'	25:RA:1376:C:H4'	2.18	0.43
25:RA:579:G:H2'	25:RA:580:C:C6	2.53	0.43
25:RA:592:G:O2'	54:R8:4:MET:CB	2.67	0.43
25:RA:593:G:C1'	54:R8:4:MET:HE1	2.48	0.43
25:RA:638:G:H2'	25:RA:639:U:C6	2.54	0.43
25:RA:666:G:H4'	35:RP:49:ARG:NH1	2.34	0.43
25:RA:1270:C:H5"	25:RA:1271:G:O5'	2.19	0.43
25:RA:1509:C:H3'	25:RA:1510:A:H5"	2.00	0.43
25:RA:1816:G:C8	27:RD:62:TYR:CZ	3.07	0.43
25:RA:2893:G:H5"	25:RA:2894:G:H5'	2.00	0.43
29:RF:101:LEU:HD12	29:RF:102:PRO:HD2	2.00	0.43
29:RF:148:LEU:HD11	29:RF:193:VAL:HG21	2.01	0.43
30:RG:67:LYS:HE2	50:R4:6:HIS:NE2	2.33	0.43
32:RI:21:VAL:HG21	32:RI:25:TYR:HD2	1.84	0.43
33:RN:58:ASP:HB3	33:RN:95:PRO:HB3	2.00	0.43
33:RN:61:ARG:HA	33:RN:61:ARG:HE	1.82	0.43
34:RO:104:ARG:HD3	39:RT:36:GLU:OE2	2.18	0.43
38:RS:12:PHE:HD1	38:RS:12:PHE:HA	1.72	0.43
1:XA:950:U:H2'	1:XA:951:G:H8	1.84	0.43
1:XA:1250:A:H2	1:XA:1370:G:H1'	1.84	0.43
3:XC:149:ALA:HA	3:XC:201:TYR:O	2.18	0.43
5:XE:42:GLY:CA	5:XE:66:MET:HG2	2.48	0.43
9:XI:8:GLY:HA2	9:XI:79:LEU:HD12	2.01	0.43
10:XJ:6:ILE:HG22	10:XJ:98:ILE:HG23	2.01	0.43
23:XX:19:A:C2'	23:XX:20:C:H5'	2.48	0.43
25:YA:265:A:O2'	25:YA:266:G:H4'	2.18	0.43
25:YA:459:U:H5"	53:Y7:40:TRP:CD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:969:U:H2'	25:YA:970:C:C6	2.53	0.43
25:YA:1101:U:H2'	25:YA:1102:C:C6	2.53	0.43
25:YA:1952:A:C6	34:YO:22:ILE:HD12	2.54	0.43
27:YD:172:TYR:HD1	27:YD:185:VAL:C	2.26	0.43
28:YE:13:ARG:HH11	28:YE:13:ARG:HB2	1.81	0.43
28:YE:23:VAL:HG12	28:YE:184:VAL:O	2.19	0.43
31:YH:119:GLU:CD	31:YH:120:GLY:N	2.77	0.43
31:YH:120:GLY:O	31:YH:136:ILE:HD12	2.19	0.43
35:YP:65:ARG:HH21	54:Y8:46:ARG:NH1	2.06	0.43
35:YP:97:PRO:HD3	35:YP:126:VAL:O	2.18	0.43
48:Y2:4:SER:OG	48:Y2:5:GLU:OE2	2.26	0.43
1:QA:105:G:C6	1:QA:106:C:C4	3.07	0.43
1:QA:757:U:O2'	1:QA:879:C:H1'	2.18	0.43
1:QA:818:G:H3'	1:QA:819:A:H5"	2.01	0.43
1:QA:1494:G:N7	58:QA:1696:PAR:N32	2.66	0.43
3:QC:11:ARG:HB3	3:QC:15:THR:HB	2.01	0.43
12:QL:76:ASN:CG	12:QL:76:ASN:O	2.61	0.43
15:QO:48:LYS:HD3	15:QO:48:LYS:HA	1.75	0.43
17:QQ:63:ARG:HG2	17:QQ:64:PRO:HD2	2.00	0.43
25:RA:27:G:H1'	25:RA:513:A:H61	1.82	0.43
25:RA:1003:G:O2'	25:RA:1010:A:N1	2.36	0.43
25:RA:1032:A:H2	25:RA:1122:G:H1	1.63	0.43
25:RA:1131:G:C8	25:RA:2025:C:H4'	2.54	0.43
25:RA:2014:A:H2'	25:RA:2015:A:C8	2.54	0.43
25:RA:2193:G:H2'	25:RA:2194:G:O4'	2.19	0.43
25:RA:2705:A:O2'	25:RA:2852:G:OP1	2.34	0.43
26:RB:65:C:C4	26:RB:108:C:C4	3.07	0.43
31:RH:35:VAL:CG2	31:RH:75:ALA:HB2	2.48	0.43
31:RH:119:GLU:CD	31:RH:120:GLY:N	2.77	0.43
32:RI:30:LEU:HD22	32:RI:35:LEU:HD11	2.00	0.43
32:RI:72:LEU:HD11	32:RI:101:LEU:HD11	2.00	0.43
35:RP:65:ARG:NE	54:R8:15:LYS:HB2	2.34	0.43
40:RU:69:CYS:HB3	40:RU:106:PHE:CZ	2.54	0.43
44:RY:54:LYS:HB3	44:RY:55:TYR:CD1	2.53	0.43
45:RZ:141:VAL:HA	45:RZ:144:LEU:HD23	2.00	0.43
1:XA:437:U:H2'	1:XA:438:G:O4'	2.18	0.43
1:XA:1144:G:N2	1:XA:1146:A:H62	2.17	0.43
1:XA:1317:C:H5"	1:XA:1318:A:OP2	2.19	0.43
1:XA:1443:G:H5'	1:XA:1446:A:OP2	2.18	0.43
2:XB:7:VAL:HG21	2:XB:217:ARG:NH1	2.34	0.43
2:XB:172:ILE:O	2:XB:175:ARG:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:XB:212:GLN:NE2	2:XB:216:SER:HB2	2.34	0.43
7:XG:140:ASP:HA	7:XG:143:ARG:NH1	2.34	0.43
8:XH:13:ILE:O	8:XH:17:THR:HG23	2.18	0.43
8:XH:104:ARG:HD2	8:XH:138:TRP:CD2	2.53	0.43
13:XM:4:ILE:HG22	13:XM:5:ALA:N	2.34	0.43
14:XN:6:LEU:HD23	14:XN:23:ARG:HH22	1.83	0.43
25:YA:34:C:N4	25:YA:447:A:H61	2.17	0.43
25:YA:222:A:H3'	25:YA:421:U:C5'	2.47	0.43
25:YA:271:G:H2'	25:YA:272:G:H8	1.83	0.43
25:YA:1359:A:H5'	25:YA:1359:A:C8	2.54	0.43
25:YA:1534:G:H2'	25:YA:1534:G:N3	2.34	0.43
25:YA:1641:A:H2'	25:YA:1642:G:O4'	2.19	0.43
25:YA:2356:C:H2'	25:YA:2357:U:O4'	2.18	0.43
25:YA:2712:U:O2'	25:YA:2712(A):A:C8	2.69	0.43
26:YB:29:A:H2'	26:YB:30:C:C6	2.54	0.43
27:YD:117:VAL:HG21	27:YD:128:GLY:C	2.44	0.43
29:YF:64:ILE:HD12	29:YF:64:ILE:HA	1.89	0.43
30:YG:31:VAL:HA	30:YG:32:PRO:HD3	1.83	0.43
30:YG:103:LEU:HA	30:YG:103:LEU:HD23	1.83	0.43
32:YI:101:LEU:HD22	32:YI:107:VAL:HB	2.00	0.43
36:YQ:57:HIS:ND1	36:YQ:58:PHE:N	2.67	0.43
37:YR:91:GLN:H	37:YR:91:GLN:CD	2.26	0.43
44:YY:80:GLY:O	44:YY:81:LYS:HG3	2.18	0.43
52:Y6:28:ARG:HH21	52:Y6:30:THR:HG23	1.84	0.43
1:QA:514:C:H2'	1:QA:515:G:C8	2.53	0.43
1:QA:895:G:H1	1:QA:904:C:N4	2.16	0.43
1:QA:983:A:H5"	1:QA:984:C:OP2	2.19	0.43
4:QD:22:LYS:HB2	4:QD:26:CYS:HB2	2.00	0.43
4:QD:78:LEU:HA	4:QD:78:LEU:HD23	1.76	0.43
8:QH:101:PRO:HG2	8:QH:133:LEU:HD11	2.01	0.43
11:QK:53:SER:C	11:QK:55:LYS:H	2.27	0.43
15:QO:25:THR:HG21	15:QO:70:LEU:HB2	2.00	0.43
17:QQ:45:HIS:NE2	17:QQ:47:PRO:HG3	2.34	0.43
25:RA:177:G:H5"	25:RA:177:G:N3	2.34	0.43
25:RA:207:A:H2'	25:RA:208:C:O4'	2.19	0.43
25:RA:601:C:O2	25:RA:605:C:H4'	2.19	0.43
25:RA:1204:A:HO2'	25:RA:1205:U:P	2.41	0.43
25:RA:2300:G:H2'	25:RA:2301:C:C6	2.54	0.43
25:RA:2582:G:N2	25:RA:2583:G:H1'	2.33	0.43
31:RH:16:SER:OG	31:RH:17:VAL:N	2.50	0.43
33:RN:89:LYS:O	33:RN:93:THR:HG22	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:RT:107:ASP:O	39:RT:110:ILE:HG22	2.19	0.43
46:R0:51:VAL:N	46:R0:62:LEU:HD12	2.33	0.43
50:R4:43:TYR:O	50:R4:46:GLN:HA	2.19	0.43
52:R6:7:ILE:HG13	52:R6:8:LYS:H	1.84	0.43
4:XD:196:LEU:O	4:XD:198:VAL:N	2.51	0.43
5:XE:79:GLU:H	5:XE:79:GLU:HG3	1.45	0.43
25:YA:900:A:H3'	25:YA:901:A:H8	1.84	0.43
25:YA:1759:A:H1'	25:YA:2711:A:C2	2.54	0.43
25:YA:1930:G:H2'	25:YA:1968:G:C6	2.54	0.43
25:YA:2832:U:O4	25:YA:2884:U:H5'	2.18	0.43
26:YB:49:C:H2'	26:YB:50:G:C8	2.53	0.43
28:YE:52:LEU:O	28:YE:74:PRO:HA	2.19	0.43
28:YE:179:GLU:CB	28:YE:181:LEU:HD23	2.24	0.43
29:YF:45:ARG:HH11	29:YF:45:ARG:HG2	1.82	0.43
29:YF:63:LYS:CE	29:YF:67:GLN:HB2	2.49	0.43
31:YH:58:GLU:O	31:YH:59:ARG:C	2.61	0.43
31:YH:125:VAL:HG12	31:YH:126:PRO:CD	2.49	0.43
36:YQ:25:ASP:H	36:YQ:102:VAL:HG23	1.84	0.43
41:YV:64:HIS:ND1	41:YV:92:THR:HG22	2.34	0.43
45:YZ:102:LEU:HG	45:YZ:123:ASP:HA	2.01	0.43
49:Y3:7:LYS:HE2	49:Y3:32:GLN:O	2.19	0.43
50:Y4:6:HIS:HA	50:Y4:7:PRO:HD2	1.82	0.43
1:QA:584:G:H2'	1:QA:585:G:H8	1.84	0.42
1:QA:1095:U:H5"	1:QA:1109:C:O2	2.18	0.42
3:QC:148:GLY:HA3	3:QC:172:ARG:O	2.18	0.42
5:QE:18:ARG:HE	5:QE:18:ARG:HB3	1.52	0.42
6:QF:62:TRP:CH2	6:QF:64:GLN:HB2	2.54	0.42
9:QI:17:VAL:HG11	9:QI:81:ILE:HD13	2.00	0.42
14:QN:47:LEU:HD23	14:QN:47:LEU:HA	1.74	0.42
22:QV:76:A:O2'	56:Z5:101:PPU:O	2.26	0.42
25:RA:709:U:H2'	25:RA:710:G:C8	2.54	0.42
25:RA:1445:C:H2'	25:RA:1446:C:C6	2.54	0.42
25:RA:2298:A:N6	25:RA:2318:G:H8	2.15	0.42
25:RA:2651:C:N4	25:RA:2669:G:H1	2.13	0.42
27:RD:28:GLU:H	27:RD:28:GLU:CD	2.27	0.42
27:RD:169:GLU:N	27:RD:172:TYR:O	2.51	0.42
28:RE:23:VAL:HG12	28:RE:184:VAL:O	2.19	0.42
37:RR:113:LEU:HD12	37:RR:113:LEU:HA	1.88	0.42
40:RU:75:ASN:HB2	40:RU:78:THR:H	1.84	0.42
50:R4:2:LYS:HD2	50:R4:2:LYS:HA	1.61	0.42
50:R4:48:ARG:NH1	50:R4:51:ASP:HA	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:R8:28:GLY:O	54:R8:29:LYS:C	2.62	0.42
1:XA:511:C:O2'	1:XA:534:U:H1'	2.19	0.42
1:XA:515:G:N2	1:XA:537:G:C4	2.87	0.42
1:XA:731:G:OP1	1:XA:766:A:H1'	2.19	0.42
1:XA:1437:C:N3	1:XA:1464:G:N2	2.61	0.42
2:XB:19:HIS:HD2	2:XB:20:GLU:CD	2.27	0.42
2:XB:98:LEU:HG	2:XB:101:MET:HE3	2.01	0.42
18:XR:56:THR:HB	18:XR:58:LEU:HD12	2.01	0.42
20:XT:33:ILE:CG2	20:XT:63:ILE:HG12	2.49	0.42
25:YA:20:C:H42	25:YA:520:G:H1	1.66	0.42
25:YA:570:G:H2'	25:YA:2030:A:C5	2.54	0.42
25:YA:1448:G:H1'	25:YA:1528:A:H62	1.83	0.42
25:YA:2318:G:N2	38:YS:2:ALA:HA	2.34	0.42
25:YA:2712:U:H1'	25:YA:2712(A):A:C8	2.53	0.42
27:YD:17:THR:HG22	27:YD:204:ILE:HA	1.98	0.42
27:YD:33:LEU:O	27:YD:35:LYS:N	2.52	0.42
29:YF:135:LYS:O	29:YF:136:THR:C	2.61	0.42
35:YP:96:THR:O	35:YP:99:LEU:HB3	2.18	0.42
36:YQ:83:MET:HB2	46:Y0:7:LEU:HB3	2.01	0.42
45:YZ:20:ARG:O	45:YZ:20:ARG:HD3	2.18	0.42
47:Y1:58:ILE:HG23	47:Y1:87:PRO:HG3	2.01	0.42
48:Y2:48:HIS:C	48:Y2:50:ILE:N	2.76	0.42
52:Y6:41:PRO:HG2	52:Y6:45:LYS:N	2.29	0.42
54:Y8:53:PRO:HD2	54:Y8:54:GLU:H	1.84	0.42
1:QA:694:A:N1	1:QA:787:A:O2'	2.52	0.42
1:QA:1216:G:H2'	1:QA:1217:C:C6	2.53	0.42
1:QA:1319:A:OP1	19:QS:3:ARG:HD3	2.19	0.42
1:QA:1321:C:H5'	13:QM:87:TYR:CE1	2.54	0.42
1:QA:1442:G:C5	1:QA:1446:A:C6	3.07	0.42
1:QA:1443:G:H5'	1:QA:1446:A:OP2	2.19	0.42
7:QG:102:ARG:HG2	7:QG:106:GLN:OE1	2.20	0.42
10:QJ:76:ASN:HA	10:QJ:77:PRO:HD2	1.85	0.42
19:QS:41:VAL:HG12	19:QS:44:MET:HB2	2.01	0.42
20:QT:87:LYS:HD2	20:QT:87:LYS:HA	1.68	0.42
25:RA:223:A:N1	25:RA:407:G:O2'	2.43	0.42
25:RA:270(R):G:H2'	25:RA:270(S):G:C8	2.54	0.42
25:RA:654(A):G:H1	25:RA:654(T):C:N4	2.17	0.42
25:RA:1432:C:H2'	25:RA:1433:U:O4'	2.19	0.42
25:RA:1482:U:H5'	25:RA:1483:G:OP2	2.18	0.42
25:RA:1709:U:H2'	25:RA:1710:C:C6	2.54	0.42
25:RA:2455:G:H2'	25:RA:2456:C:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2466:C:OP1	55:R9:4:ARG:HB2	2.18	0.42
26:RB:44:G:H5"	26:RB:45:A:OP1	2.19	0.42
26:RB:74:U:H2'	26:RB:75:G:O4'	2.19	0.42
28:RE:155:LYS:O	28:RE:156:MET:HG3	2.19	0.42
30:RG:97:ASP:HA	30:RG:100:TRP:HD1	1.84	0.42
33:RN:30:ILE:HG23	33:RN:52:VAL:HG11	1.99	0.42
34:RO:104:ARG:HH22	39:RT:43:GLN:CD	2.26	0.42
35:RP:61:ARG:HG3	54:R8:13:ARG:CD	2.49	0.42
38:RS:93:LYS:HE3	38:RS:93:LYS:HB2	1.65	0.42
51:R5:3:LYS:HA	51:R5:3:LYS:CE	2.36	0.42
51:R5:35:GLU:O	51:R5:36:CYS:C	2.63	0.42
52:R6:8:LYS:O	52:R6:27:LYS:HA	2.18	0.42
52:R6:35:GLU:H	52:R6:35:GLU:HG2	1.72	0.42
1:XA:106:C:O2	1:XA:379:C:H4'	2.19	0.42
1:XA:619:U:H3	4:XD:135:LEU:CD2	2.25	0.42
1:XA:912:C:O2'	1:XA:913:A:H5'	2.18	0.42
2:XB:9:GLU:H	2:XB:9:GLU:CD	2.27	0.42
6:XF:30:LEU:HB3	6:XF:35:ALA:HB3	2.01	0.42
25:YA:139:G:H1'	25:YA:140:A:H2	1.84	0.42
25:YA:610:C:H42	25:YA:618:G:H1	1.65	0.42
25:YA:686:G:N2	25:YA:788:A:H61	2.17	0.42
25:YA:1438:U:H2'	25:YA:1439:A:H8	1.82	0.42
25:YA:1467:C:C5	25:YA:1546:C:H2'	2.53	0.42
25:YA:1729:A:N6	25:YA:1731:G:N3	2.67	0.42
25:YA:1844:C:H2'	25:YA:1845:G:C8	2.53	0.42
25:YA:2453:A:H2'	25:YA:2454:G:H8	1.83	0.42
30:YG:99:MET:HB2	30:YG:99:MET:HE3	1.66	0.42
38:YS:64:GLU:O	38:YS:68:GLN:HG3	2.19	0.42
45:YZ:74:VAL:HG13	45:YZ:86:VAL:HG22	2.00	0.42
45:YZ:150:LEU:HB2	45:YZ:154:ASP:OD2	2.19	0.42
50:Y4:14:ILE:HG13	50:Y4:31:ILE:HB	1.99	0.42
53:Y7:25:PRO:HA	53:Y7:28:ARG:CZ	2.49	0.42
1:QA:1163:C:H5'	1:QA:1164:G:OP2	2.20	0.42
1:QA:1402:C:O2	1:QA:1500:A:N1	2.53	0.42
2:QB:21:ARG:HG3	2:QB:38:GLY:O	2.20	0.42
3:QC:59:ARG:HH12	3:QC:97:LYS:HE3	1.84	0.42
3:QC:112:SER:O	3:QC:116:VAL:HG23	2.20	0.42
8:QH:105:ARG:HA	8:QH:105:ARG:HD3	1.78	0.42
9:QI:48:GLU:N	9:QI:49:PRO:HD2	2.35	0.42
17:QQ:29:HIS:CG	17:QQ:30:PRO:HD2	2.54	0.42
25:RA:856:C:N4	25:RA:921:G:H1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:2643:G:H2'	25:RA:2644:G:O4'	2.19	0.42
27:RD:245:PRO:HA	27:RD:246:PRO:HD3	1.95	0.42
28:RE:3:GLY:HA3	28:RE:81:ILE:CD1	2.47	0.42
29:RF:51:THR:HG23	29:RF:92:PRO:HG2	2.02	0.42
29:RF:107:LYS:HB3	29:RF:107:LYS:HE2	1.77	0.42
31:RH:26:VAL:CG1	31:RH:33:LEU:HB2	2.50	0.42
31:RH:58:GLU:O	31:RH:60:ARG:N	2.53	0.42
31:RH:120:GLY:O	31:RH:136:ILE:HD12	2.19	0.42
32:RI:135:GLU:HB2	32:RI:136:VAL:H	1.60	0.42
33:RN:96:GLU:HB2	33:RN:122:VAL:HG12	2.00	0.42
36:RQ:65:PHE:O	36:RQ:66:ILE:CG1	2.48	0.42
41:RV:64:HIS:CG	41:RV:92:THR:HG22	2.52	0.42
45:RZ:52:SER:O	45:RZ:52:SER:OG	2.33	0.42
48:R2:70:GLN:O	48:R2:71:ASN:HB2	2.19	0.42
52:R6:17:LYS:HB3	52:R6:44:ARG:NH2	2.30	0.42
1:XA:457:C:H2'	1:XA:458:C:C6	2.54	0.42
1:XA:563:A:N7	1:XA:567:G:H1'	2.34	0.42
1:XA:987:G:H1	1:XA:1218:C:N4	2.15	0.42
2:XB:68:ILE:HB	2:XB:70:PHE:HE1	1.84	0.42
4:XD:112:VAL:N	4:XD:116:GLN:OE1	2.38	0.42
5:XE:33:VAL:O	5:XE:112:LEU:HD12	2.20	0.42
7:XG:38:LEU:HD12	7:XG:38:LEU:O	2.20	0.42
13:XM:117:VAL:O	13:XM:118:ALA:C	2.62	0.42
20:XT:44:ALA:O	20:XT:91:LEU:O	2.37	0.42
25:YA:270(J):G:H2'	25:YA:270(K):C:O4'	2.19	0.42
25:YA:279:C:H42	25:YA:361:G:H1	1.67	0.42
25:YA:896:A:C4	45:YZ:146:ILE:HD12	2.54	0.42
25:YA:1109:C:O2'	25:YA:1110:G:OP1	2.33	0.42
25:YA:1728:G:H3'	25:YA:1729:A:C5'	2.46	0.42
26:YB:16:G:H1	26:YB:68:C:H42	1.67	0.42
27:YD:108:PRO:HG2	27:YD:111:LEU:HB2	2.01	0.42
30:YG:16:ARG:N	30:YG:17:PRO:HD2	2.34	0.42
30:YG:64:THR:CG2	30:YG:66:GLN:H	2.28	0.42
31:YH:125:VAL:CG1	31:YH:126:PRO:CG	2.94	0.42
32:YI:24:GLY:O	32:YI:28:ASN:HB2	2.19	0.42
39:YT:80:SER:HA	39:YT:81:PRO:HD3	1.89	0.42
50:Y4:60:GLN:O	50:Y4:63:TYR:HB3	2.20	0.42
52:Y6:10:LEU:HG	54:Y8:34:TRP:CD1	2.54	0.42
52:Y6:15:GLU:HG2	52:Y6:49:HIS:NE2	2.34	0.42
1:QA:27:G:H4'	4:QD:209:ARG:HG3	2.00	0.42
1:QA:181:G:O2'	1:QA:182:U:O5'	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:401:C:H2'	1:QA:402:G:H8	1.83	0.42
1:QA:519:C:H2'	1:QA:520:A:O4'	2.19	0.42
1:QA:1479:C:H2'	1:QA:1480:G:H8	1.84	0.42
2:QB:44:LEU:H	2:QB:44:LEU:HD12	1.83	0.42
2:QB:69:LEU:O	2:QB:162:ILE:HA	2.19	0.42
14:QN:41:ARG:NH2	14:QN:42:ILE:HD11	2.35	0.42
19:QS:10:PHE:CG	19:QS:11:VAL:N	2.88	0.42
25:RA:297:C:H5"	44:RY:85:VAL:CG2	2.49	0.42
25:RA:1636:C:H2'	25:RA:1637:A:H8	1.82	0.42
25:RA:2116:G:H2'	25:RA:2117:A:C8	2.54	0.42
25:RA:2682:U:O2'	28:RE:13:ARG:CG	2.65	0.42
25:RA:2683:C:O2	34:RO:70:LYS:NZ	2.39	0.42
25:RA:2845:G:H5"	39:RT:55:ASN:HA	2.01	0.42
28:RE:28:ALA:HB3	28:RE:93:VAL:CG2	2.46	0.42
28:RE:72:VAL:O	28:RE:73:GLU:C	2.62	0.42
28:RE:76:ARG:C	28:RE:78:LEU:N	2.76	0.42
28:RE:104:VAL:CG1	28:RE:188:VAL:HG23	2.49	0.42
28:RE:137:HIS:CB	28:RE:138:PRO:HD2	2.42	0.42
28:RE:188:VAL:HA	28:RE:189:PRO:HD2	1.79	0.42
32:RI:88:ILE:HG12	32:RI:88:ILE:H	1.70	0.42
50:R4:68:ARG:O	50:R4:69:LYS:HB2	2.17	0.42
51:R5:20:ARG:HA	51:R5:23:HIS:CE1	2.54	0.42
52:R6:45:LYS:HD3	52:R6:45:LYS:HA	1.75	0.42
1:XA:255:G:H4'	17:XQ:17:LYS:HD3	2.00	0.42
1:XA:660:G:C2	1:XA:746:A:C2	3.08	0.42
1:XA:765:G:N2	1:XA:813:U:OP2	2.52	0.42
1:XA:1129:C:O2	1:XA:1132:C:N4	2.38	0.42
1:XA:1483:A:HO2'	25:YA:1947:C:HO2'	1.61	0.42
1:XA:1504:G:H4'	1:XA:1505:G:C4	2.54	0.42
4:XD:16:GLY:HA2	4:XD:33:MET:HE1	2.00	0.42
6:XF:95:GLU:HA	6:XF:96:PRO:HD3	1.87	0.42
22:XV:15:C:H5'	22:XV:16:C:H5	1.84	0.42
25:YA:833:U:H2'	25:YA:834:C:C6	2.53	0.42
25:YA:1680:U:O2'	25:YA:1763:G:N7	2.44	0.42
25:YA:2360:A:H2'	25:YA:2361:A:O4'	2.20	0.42
25:YA:2572:A:N3	28:YE:144:ARG:NH2	2.68	0.42
26:YB:24:G:O6	26:YB:56:G:O2'	2.30	0.42
28:YE:36:ARG:HH11	28:YE:36:ARG:CB	2.28	0.42
28:YE:70:ALA:O	28:YE:71:GLY:C	2.62	0.42
28:YE:104:VAL:CG1	28:YE:188:VAL:HG23	2.49	0.42
28:YE:117:MET:HA	28:YE:122:PHE:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:197:ILE:CD1	28:YE:199:ARG:HH12	2.26	0.42
29:YF:32:LEU:C	29:YF:32:LEU:CD1	2.83	0.42
29:YF:128:ALA:O	29:YF:129:PHE:CB	2.67	0.42
33:YN:96:GLU:HG2	33:YN:97:ARG:H	1.84	0.42
36:YQ:20:ALA:HA	36:YQ:98:LYS:HB3	2.02	0.42
36:YQ:20:ALA:C	36:YQ:21:THR:HG23	2.43	0.42
38:YS:15:ARG:O	38:YS:19:LYS:HD3	2.20	0.42
38:YS:110:LEU:HD23	38:YS:112:PHE:CE1	2.55	0.42
39:YT:45:PHE:CE1	39:YT:65:LYS:HE3	2.55	0.42
42:YW:86:LEU:HD22	42:YW:96:ILE:HD12	2.01	0.42
43:YX:87:GLN:C	43:YX:88:LYS:HG3	2.42	0.42
46:Y0:53:MET:CB	46:Y0:59:LEU:HD23	2.50	0.42
48:Y2:27:GLU:N	48:Y2:27:GLU:CD	2.77	0.42
49:Y3:4:LEU:HD22	49:Y3:56:VAL:HG12	2.01	0.42
1:QA:129(A):G:N2	1:QA:188:U:O2'	2.52	0.42
1:QA:585:G:O3'	17:QQ:34:LYS:NZ	2.51	0.42
1:QA:827:U:C5	1:QA:870:U:C2	3.07	0.42
1:QA:960:U:H4'	1:QA:961:U:C5'	2.50	0.42
1:QA:984:C:H2'	1:QA:985:C:C6	2.55	0.42
1:QA:1040:U:H2'	1:QA:1041:A:O4'	2.19	0.42
1:QA:1109:C:OP2	3:QC:176:HIS:ND1	2.51	0.42
1:QA:1163:C:N4	1:QA:1173:G:H1	2.17	0.42
1:QA:1516:G:H2'	1:QA:1518:A:OP2	2.19	0.42
2:QB:92:TYR:CD1	2:QB:151:GLY:HA3	2.55	0.42
17:QQ:10:VAL:HG13	17:QQ:19:VAL:HB	2.01	0.42
25:RA:95:G:H1'	48:R2:47:ASN:HB3	2.00	0.42
25:RA:569:U:O2'	25:RA:983:A:N1	2.37	0.42
25:RA:594:U:H5'	54:R8:61:LEU:HD21	2.00	0.42
25:RA:702:G:C2	25:RA:731:C:C2	3.07	0.42
25:RA:741:G:H2'	25:RA:742:G:C8	2.54	0.42
25:RA:1210:A:H4'	25:RA:1211:U:O5'	2.19	0.42
25:RA:2074:U:H2'	25:RA:2075:U:C6	2.55	0.42
25:RA:2086:U:H2'	25:RA:2087:G:C8	2.55	0.42
25:RA:2242:G:H2'	25:RA:2243:U:O4'	2.19	0.42
25:RA:2331:G:O2'	46:R0:43:THR:HG22	2.19	0.42
25:RA:2784:C:H5"	28:RE:41:LYS:NZ	2.34	0.42
27:RD:96:HIS:NE2	27:RD:102:LYS:HE2	2.34	0.42
28:RE:25:VAL:HG11	39:RT:11:GLU:CD	2.44	0.42
28:RE:35:GLN:HB3	28:RE:48:GLN:HB2	2.01	0.42
28:RE:144:ARG:HB3	28:RE:145:LYS:H	1.58	0.42
30:RG:159:VAL:HG21	30:RG:173:LEU:HD11	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:RI:82:ARG:HD3	32:RI:146:ALA:HB3	2.01	0.42
33:RN:56:ASN:C	33:RN:58:ASP:H	2.27	0.42
40:RU:83:LEU:HD12	40:RU:113:ALA:HB2	2.01	0.42
45:RZ:93:ASP:OD1	45:RZ:93:ASP:N	2.53	0.42
50:R4:35:VAL:C	50:R4:37:SER:N	2.77	0.42
50:R4:54:GLY:HA2	50:R4:57:GLU:CG	2.50	0.42
52:R6:11:LEU:HD13	52:R6:11:LEU:HA	1.81	0.42
1:XA:254:G:H2'	1:XA:255:G:H8	1.84	0.42
1:XA:665:A:H1'	1:XA:733:A:O4'	2.19	0.42
1:XA:965:A:H4'	1:XA:966:G:OP1	2.19	0.42
1:XA:1029:G:O2'	1:XA:1032(A):G:N1	2.43	0.42
2:XB:212:GLN:CD	2:XB:235:SER:HB2	2.44	0.42
3:XC:85:ARG:HD2	3:XC:85:ARG:HA	1.83	0.42
9:XI:32:ASP:OD1	9:XI:33:PHE:N	2.53	0.42
13:XM:7:VAL:O	13:XM:9:ILE:HG23	2.19	0.42
13:XM:40:ASN:ND2	13:XM:43:THR:HG23	2.34	0.42
20:XT:43:LEU:HD23	20:XT:43:LEU:HA	1.88	0.42
25:YA:31:C:O2'	25:YA:1238:G:H5'	2.19	0.42
25:YA:1153:C:H5'	40:YU:76:TYR:HE1	1.85	0.42
25:YA:1188:U:O2'	25:YA:1189:A:H5'	2.19	0.42
25:YA:1454:U:H5'	37:YR:63:ARG:NE	2.35	0.42
27:YD:71:ASP:CB	27:YD:103:ARG:HH22	2.32	0.42
27:YD:155:LEU:HD23	27:YD:177:LEU:HD21	2.00	0.42
27:YD:215:LEU:H	27:YD:215:LEU:HG	1.59	0.42
28:YE:28:ALA:HB3	28:YE:93:VAL:CG2	2.47	0.42
31:YH:58:GLU:O	31:YH:60:ARG:N	2.52	0.42
31:YH:77:LYS:HB3	31:YH:77:LYS:HZ2	1.78	0.42
32:YI:14:ASP:OD1	32:YI:14:ASP:N	2.52	0.42
33:YN:134:ARG:H	33:YN:135:PRO:HD3	1.83	0.42
38:YS:42:ASP:C	38:YS:44:LYS:N	2.72	0.42
50:Y4:38:LYS:C	50:Y4:40:HIS:H	2.23	0.42
51:Y5:58:LEU:HD13	51:Y5:60:VAL:OXT	2.19	0.42
1:QA:429:U:H1'	1:QA:430:A:H5"	2.02	0.42
1:QA:878:G:H5'	8:QH:89:PRO:HG2	2.02	0.42
1:QA:1098:C:H2'	1:QA:1099:G:O4'	2.19	0.42
1:QA:1226:C:H2'	13:QM:103:THR:HB	2.02	0.42
1:QA:1312:G:OP2	50:R4:67:TYR:HE1	2.01	0.42
1:QA:1346:A:C4	7:QG:10:ARG:NH1	2.88	0.42
2:QB:60:ASP:O	2:QB:64:ARG:HG2	2.19	0.42
6:QF:22:GLU:O	6:QF:26:ILE:HG13	2.19	0.42
6:QF:99:ALA:HB1	18:QR:23:LYS:HZ2	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:QL:38:THR:HG22	12:QL:57:LYS:HB3	2.01	0.42
13:QM:4:ILE:H	13:QM:9:ILE:HG22	1.84	0.42
25:RA:270(I):G:H2'	25:RA:270(J):G:H8	1.84	0.42
25:RA:503:A:H4'	25:RA:504:U:H5"	2.01	0.42
25:RA:1007:C:H5"	33:RN:35:ARG:HH11	1.84	0.42
25:RA:1046:A:N3	25:RA:1046:A:H3'	2.34	0.42
25:RA:1092:C:O2'	31:RH:170:ARG:CG	2.68	0.42
25:RA:1329:U:H5"	25:RA:1330:C:H5	1.83	0.42
25:RA:1939:U:O4'	25:RA:2591:C:O2'	2.37	0.42
25:RA:2224:G:OP1	27:RD:268:ARG:HD3	2.19	0.42
27:RD:33:LEU:HB3	27:RD:34:VAL:H	1.64	0.42
27:RD:123:ALA:HA	27:RD:124:PRO:HD2	1.77	0.42
28:RE:24:THR:HB	28:RE:184:VAL:HG23	2.02	0.42
29:RF:28:ILE:H	29:RF:28:ILE:HG13	1.69	0.42
31:RH:89:ILE:CD1	31:RH:89:ILE:H	2.32	0.42
31:RH:125:VAL:HG12	31:RH:126:PRO:CD	2.49	0.42
31:RH:136:ILE:O	31:RH:137:ASP:O	2.38	0.42
32:RI:93:THR:O	32:RI:97:ILE:HG12	2.19	0.42
33:RN:17:ASP:C	33:RN:19:GLU:H	2.28	0.42
33:RN:35:ARG:HB2	33:RN:42:TRP:CZ3	2.54	0.42
35:RP:6:LEU:HA	35:RP:6:LEU:HD13	1.75	0.42
36:RQ:27:VAL:HG11	36:RQ:134:ARG:HG3	2.01	0.42
36:RQ:139:GLU:H	36:RQ:139:GLU:CD	2.28	0.42
38:RS:39:ILE:HD11	38:RS:73:LEU:HD11	2.00	0.42
44:RY:50:ARG:H	44:RY:50:ARG:HG2	1.67	0.42
1:XA:61:G:H2'	1:XA:62:U:O4'	2.19	0.42
1:XA:601:C:H2'	1:XA:602:A:C8	2.54	0.42
1:XA:990:C:H2'	1:XA:991:U:C6	2.54	0.42
1:XA:1446:A:HO2'	1:XA:1447:G:P	2.43	0.42
2:XB:113:HIS:O	2:XB:116:GLU:HB2	2.20	0.42
3:XC:32:LEU:HD22	3:XC:59:ARG:NH1	2.34	0.42
4:XD:120:LEU:HD23	4:XD:120:LEU:HA	1.89	0.42
18:XR:43:PHE:CE1	18:XR:58:LEU:HD11	2.54	0.42
22:XV:12:C:O2'	25:YA:1924:C:H4'	2.19	0.42
25:YA:184:C:H2'	25:YA:185:U:H6	1.84	0.42
25:YA:724:U:H2'	25:YA:725:G:O4'	2.19	0.42
25:YA:900:A:H3'	25:YA:901:A:C8	2.54	0.42
25:YA:1210:A:H5"	25:YA:1212:G:O4'	2.19	0.42
25:YA:1342:A:OP1	43:YX:36:LYS:NZ	2.52	0.42
25:YA:2453:A:H2'	25:YA:2454:G:C8	2.55	0.42
25:YA:2500:U:O2'	25:YA:2504:U:OP1	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:2:ALA:O	27:YD:3:VAL:CB	2.68	0.42
27:YD:158:ALA:HB3	27:YD:161:THR:CG2	2.49	0.42
27:YD:263:ARG:NH1	27:YD:263:ARG:CB	2.75	0.42
28:YE:31:CYS:HB3	28:YE:49:LEU:HG	2.01	0.42
28:YE:176:ILE:HD12	28:YE:176:ILE:N	2.35	0.42
29:YF:62:ARG:NH1	29:YF:62:ARG:CB	2.82	0.42
29:YF:123:LEU:HD12	29:YF:124:LEU:H	1.82	0.42
29:YF:183:VAL:HG22	29:YF:184:TYR:N	2.35	0.42
31:YH:26:VAL:CG1	31:YH:33:LEU:HB2	2.50	0.42
31:YH:84:SER:OG	31:YH:85:LYS:N	2.51	0.42
31:YH:89:ILE:CD1	31:YH:89:ILE:H	2.32	0.42
36:YQ:20:ALA:HB2	36:YQ:99:PRO:HD2	1.99	0.42
36:YQ:27:VAL:HG11	36:YQ:134:ARG:HG3	2.00	0.42
36:YQ:51:ARG:O	36:YQ:52:VAL:C	2.62	0.42
38:YS:18:ILE:O	38:YS:19:LYS:C	2.62	0.42
52:Y6:14:THR:HG21	52:Y6:19:ARG:HH21	1.85	0.42
54:Y8:28:GLY:O	54:Y8:29:LYS:O	2.37	0.42
1:QA:13:U:C5	1:QA:20:U:O4	2.72	0.42
1:QA:397:A:H3'	1:QA:397:A:N3	2.34	0.42
1:QA:1123:A:H4'	10:QJ:36:GLY:HA3	2.02	0.42
2:QB:230:VAL:HB	2:QB:231:GLU:H	1.60	0.42
7:QG:70:LYS:HA	7:QG:71:PRO:HD2	1.89	0.42
8:QH:25:ASP:N	8:QH:25:ASP:OD1	2.50	0.42
11:QK:48:ILE:HG23	11:QK:63:LEU:HD22	2.01	0.42
15:QO:43:LEU:HA	15:QO:43:LEU:HD23	1.74	0.42
25:RA:593:G:O2'	54:R8:61:LEU:CD1	2.64	0.42
25:RA:964:C:O2'	25:RA:2273:A:N3	2.50	0.42
25:RA:997:G:OP1	40:RU:93:LYS:HB2	2.20	0.42
25:RA:1257:C:H4'	29:RF:83:PHE:CE1	2.55	0.42
25:RA:1464:C:O2'	25:RA:1528:A:H8	2.03	0.42
25:RA:2260:C:H42	25:RA:2280:G:H1	1.68	0.42
25:RA:2344:U:OP1	52:R6:38:LYS:HD3	2.20	0.42
28:RE:117:MET:HA	28:RE:122:PHE:N	2.35	0.42
30:RG:151:ALA:HB3	30:RG:153:ARG:NH1	2.35	0.42
31:RH:169:VAL:HG22	31:RH:170:ARG:N	2.26	0.42
32:RI:8:PRO:HG3	32:RI:14:ASP:HB2	2.02	0.42
35:RP:57:THR:C	35:RP:59:LEU:H	2.27	0.42
35:RP:62:LEU:HD22	35:RP:62:LEU:H	1.85	0.42
37:RR:63:ARG:HA	37:RR:80:PHE:CZ	2.54	0.42
38:RS:14:VAL:HG11	38:RS:90:GLY:O	2.19	0.42
39:RT:26:ASP:HB3	39:RT:92:GLY:N	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:RW:75:TYR:CZ	42:RW:104:THR:HG21	2.54	0.42
43:RX:44:GLU:O	43:RX:48:LYS:N	2.52	0.42
45:RZ:5:LEU:HB3	45:RZ:6:LYS:H	1.56	0.42
48:R2:35:LEU:HD11	48:R2:49:LYS:HB3	2.01	0.42
1:XA:342:C:C2	1:XA:348:G:C2	3.07	0.42
1:XA:1128:C:H4'	9:XI:16:ARG:HH12	1.85	0.42
2:XB:7:VAL:HG11	2:XB:217:ARG:CZ	2.49	0.42
2:XB:19:HIS:CE1	2:XB:206:ASP:HB2	2.54	0.42
2:XB:223:ILE:HA	2:XB:226:ARG:HB3	2.02	0.42
3:XC:91:LEU:O	3:XC:95:THR:OG1	2.19	0.42
8:XH:121:ASP:OD1	8:XH:121:ASP:N	2.46	0.42
9:XI:125:TYR:HD1	9:XI:126:SER:H	1.67	0.42
25:YA:443:A:H3'	29:YF:45:ARG:NH1	2.35	0.42
25:YA:650:C:OP1	54:Y8:48:PHE:CE1	2.72	0.42
25:YA:1419:A:C6	25:YA:1421:G:C4	3.07	0.42
25:YA:1680:U:O2	25:YA:1763:G:H3'	2.19	0.42
27:YD:12:SER:O	27:YD:14:ARG:N	2.51	0.42
27:YD:25:THR:HG23	27:YD:27:THR:HB	2.02	0.42
28:YE:35:GLN:HB3	28:YE:48:GLN:HB2	2.01	0.42
28:YE:128:SER:O	28:YE:129:HIS:HB2	2.19	0.42
28:YE:143:ASN:HB2	28:YE:147:PRO:HD2	2.00	0.42
29:YF:11:VAL:HG12	29:YF:12:LEU:H	1.85	0.42
31:YH:136:ILE:O	31:YH:137:ASP:O	2.38	0.42
34:YO:21:CYS:O	34:YO:22:ILE:HD13	2.20	0.42
38:YS:49:VAL:HG21	38:YS:77:ALA:HA	2.02	0.42
39:YT:26:ASP:HB2	39:YT:91:ARG:HA	2.00	0.42
40:YU:109:LEU:HD23	40:YU:109:LEU:HA	1.89	0.42
41:YV:65:GLY:O	41:YV:90:PRO:HA	2.20	0.42
46:Y0:41:ARG:NE	46:Y0:41:ARG:HA	2.33	0.42
48:Y2:16:LEU:C	48:Y2:16:LEU:HD12	2.45	0.42
1:QA:538:G:OP1	12:QL:113:ARG:HD2	2.20	0.42
1:QA:660:G:H1	1:QA:745:C:N4	2.17	0.42
1:QA:731:G:OP1	1:QA:766:A:H1'	2.20	0.42
1:QA:885:G:O2'	1:QA:914:A:N1	2.49	0.42
1:QA:1279:A:OP2	10:QJ:9:ARG:NH1	2.52	0.42
1:QA:1312:G:H1	1:QA:1325:C:H42	1.67	0.42
2:QB:130:ARG:HA	2:QB:131:PRO:HD3	1.81	0.42
4:QD:12:CYS:HA	4:QD:19:LEU:HD21	2.02	0.42
4:QD:78:LEU:HB3	4:QD:93:PHE:HE1	1.84	0.42
5:QE:101:ILE:CG1	5:QE:119:LEU:HD23	2.49	0.42
11:QK:41:THR:HG22	11:QK:42:TRP:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:QL:61:THR:C	12:QL:63:GLY:H	2.27	0.42
15:QO:87:ILE:HG22	15:QO:88:ARG:N	2.35	0.42
19:QS:64:GLU:HG3	50:R4:55:ARG:HH12	1.84	0.42
22:QV:0:C:O2'	46:R0:6:GLY:O	2.33	0.42
25:RA:566:U:OP1	35:RP:29:LYS:HE2	2.19	0.42
25:RA:667:U:H2'	25:RA:668:G:O4'	2.20	0.42
25:RA:1006:C:O2'	33:RN:106:MET:O	2.30	0.42
25:RA:1364:G:C8	47:R1:2:SER:N	2.88	0.42
25:RA:2308:G:C6	25:RA:2311:A:N1	2.88	0.42
25:RA:2336:A:H61	46:R0:43:THR:HG21	1.85	0.42
30:RG:65:GLY:HA2	50:R4:7:PRO:HG2	2.02	0.42
30:RG:173:LEU:O	30:RG:178:PHE:HB2	2.20	0.42
31:RH:27:LYS:HD2	31:RH:27:LYS:C	2.44	0.42
32:RI:128:LEU:N	32:RI:138:ILE:O	2.45	0.42
35:RP:65:ARG:HB2	54:R8:12:LYS:O	2.20	0.42
45:RZ:136:PHE:HE2	45:RZ:138:GLU:HB3	1.83	0.42
51:R5:40:LYS:HE2	51:R5:47:PRO:HG2	2.02	0.42
52:R6:28:ARG:HB3	52:R6:30:THR:H	1.84	0.42
1:XA:135:C:H2'	1:XA:136:C:H5'	2.01	0.42
1:XA:312:C:H2'	1:XA:313:A:H8	1.85	0.42
1:XA:1157:A:H1'	1:XA:1158:C:C4	2.55	0.42
1:XA:1223:C:P	19:XS:78:ARG:HH12	2.43	0.42
7:XG:15:ASP:O	7:XG:19:GLY:HA2	2.20	0.42
11:XK:24:SER:C	11:XK:26:ASN:H	2.28	0.42
13:XM:16:ASP:HB3	13:XM:41:PRO:HB3	2.01	0.42
19:XS:78:ARG:H	19:XS:78:ARG:HG2	1.56	0.42
25:YA:271:G:H2'	25:YA:272:G:C8	2.53	0.42
25:YA:693:C:H2'	25:YA:694:U:O4'	2.19	0.42
25:YA:1043:C:N4	25:YA:1112:G:H1	2.14	0.42
25:YA:1105:U:H2'	25:YA:1106:G:H8	1.84	0.42
25:YA:1495:A:O2'	25:YA:1579:A:H5"	2.19	0.42
25:YA:2267:A:H5"	25:YA:2268:A:H5'	2.01	0.42
28:YE:76:ARG:C	28:YE:78:LEU:N	2.76	0.42
30:YG:62:LEU:HD23	30:YG:148:MET:HE1	2.01	0.42
33:YN:137:LYS:HD2	33:YN:137:LYS:HA	1.77	0.42
35:YP:144:GLU:OE1	35:YP:144:GLU:N	2.53	0.42
36:YQ:79:LEU:HB3	46:Y0:5:LYS:NZ	2.34	0.42
38:YS:52:SER:HB2	38:YS:55:ALA:CB	2.49	0.42
38:YS:111:GLU:O	38:YS:112:PHE:HD2	2.02	0.42
40:YU:30:LYS:HA	40:YU:30:LYS:HD3	1.89	0.42
43:YX:11:PRO:HD3	48:Y2:37:PHE:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:YZ:169:GLU:HG2	45:YZ:170:THR:N	2.34	0.42
46:Y0:27:GLU:HB2	46:Y0:69:PHE:CD1	2.53	0.42
46:Y0:53:MET:HB3	46:Y0:59:LEU:HD23	2.01	0.42
49:Y3:8:LEU:HB3	49:Y3:31:LEU:HA	2.01	0.42
50:Y4:68:ARG:HB2	50:Y4:69:LYS:H	1.51	0.42
51:Y5:31:VAL:HG13	51:Y5:42:PRO:HG3	2.01	0.42
1:QA:106:C:H2'	1:QA:107:G:C8	2.54	0.42
1:QA:992:U:H4'	1:QA:993:G:O5'	2.20	0.42
1:QA:1079:G:H2'	1:QA:1080:A:C8	2.54	0.42
1:QA:1207:G:H2'	1:QA:1208:C:C6	2.55	0.42
1:QA:1305:G:H1'	1:QA:1332:A:N6	2.34	0.42
2:QB:98:LEU:HD12	2:QB:101:MET:HE3	2.01	0.42
2:QB:195:ASP:O	8:QH:68:ARG:NH2	2.53	0.42
3:QC:36:ASP:HA	3:QC:39:ILE:HD12	2.02	0.42
8:QH:51:VAL:HG21	8:QH:60:ARG:HG2	2.02	0.42
8:QH:54:ASP:O	8:QH:56:LYS:HG3	2.20	0.42
15:QO:2:PRO:HB2	15:QO:3:ILE:H	1.56	0.42
19:QS:68:GLY:O	50:R4:68:ARG:HG2	2.20	0.42
25:RA:684:G:C2	25:RA:774:A:C2	3.08	0.42
25:RA:747:U:C5	51:R5:3:LYS:HB2	2.54	0.42
25:RA:1803:A:H4'	27:RD:259:THR:CG2	2.50	0.42
25:RA:2030:A:H4'	25:RA:2031:A:H8	1.85	0.42
25:RA:2138:C:H2'	25:RA:2139:C:C6	2.55	0.42
25:RA:2674:G:H2'	25:RA:2675:A:O4'	2.20	0.42
28:RE:7:VAL:CG2	28:RE:8:LYS:H	2.11	0.42
28:RE:128:SER:O	28:RE:129:HIS:HB2	2.20	0.42
28:RE:143:ASN:HB2	28:RE:147:PRO:HD2	2.01	0.42
28:RE:176:ILE:N	28:RE:176:ILE:HD12	2.35	0.42
29:RF:110:LEU:HD11	29:RF:181:LEU:HD12	2.01	0.42
30:RG:116:ASP:OD1	30:RG:116:ASP:N	2.53	0.42
32:RI:4:ILE:HG12	32:RI:18:VAL:HG22	2.01	0.42
35:RP:97:PRO:HD3	35:RP:126:VAL:O	2.20	0.42
37:RR:33:ARG:NH2	51:R5:55:ARG:CB	2.82	0.42
39:RT:51:ARG:HG3	39:RT:98:LYS:HG3	2.02	0.42
47:R1:90:ILE:O	47:R1:94:LEU:HB2	2.20	0.42
1:XA:116:A:H61	1:XA:313:A:H1'	1.84	0.42
1:XA:256:U:H2'	1:XA:257:G:C8	2.55	0.42
1:XA:523:A:C2	1:XA:527:G:C6	3.07	0.42
1:XA:1061:G:N7	3:XC:2:GLY:HA3	2.35	0.42
11:XK:109:VAL:HG11	18:XR:84:LYS:HD3	2.02	0.42
12:XL:53:ARG:HH12	12:XL:92:ASP:CB	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:XM:3:ARG:HG3	13:XM:9:ILE:HG21	2.01	0.42
13:XM:20:THR:O	13:XM:22:ILE:N	2.51	0.42
25:YA:71:A:H5"	25:YA:72:U:H3'	2.02	0.42
25:YA:256:A:H2'	25:YA:257:A:C8	2.55	0.42
25:YA:521:G:H2'	25:YA:522:G:C8	2.52	0.42
25:YA:710:G:H2'	25:YA:711:G:H8	1.85	0.42
25:YA:857:C:H1'	46:Y0:26:TYR:CE1	2.54	0.42
25:YA:1103:A:H5'	25:YA:1104:C:C5	2.55	0.42
25:YA:1130:U:HO2'	25:YA:1131:G:P	2.43	0.42
25:YA:1230:C:H2'	25:YA:1231:G:C8	2.55	0.42
25:YA:1301:A:C8	25:YA:1303:G:C8	3.08	0.42
28:YE:72:VAL:O	28:YE:73:GLU:C	2.63	0.42
29:YF:105:VAL:O	29:YF:106:ARG:C	2.63	0.42
29:YF:132:VAL:HG23	29:YF:133:ASN:H	1.83	0.42
33:YN:46:VAL:HG13	33:YN:48:MET:HG3	2.02	0.42
35:YP:97:PRO:O	35:YP:98:GLU:HB3	2.19	0.42
43:YX:84:ALA:HB1	43:YX:85:PRO:HD2	2.02	0.42
44:YY:96:ILE:HG13	44:YY:98:VAL:H	1.85	0.42
50:Y4:37:SER:HB3	50:Y4:42:PHE:HB3	2.00	0.42
51:Y5:56:LYS:H	51:Y5:56:LYS:CD	2.29	0.42
1:QA:564:C:H5'	12:QL:10:LEU:HD13	2.02	0.42
1:QA:779:C:H2'	1:QA:780:A:O4'	2.20	0.42
1:QA:1210:C:O2'	1:QA:1213:A:O2'	2.38	0.42
2:QB:27:LYS:HD2	2:QB:193:ASP:CB	2.46	0.42
2:QB:178:ARG:NH2	8:QH:74:PRO:HG3	2.34	0.42
3:QC:56:ASP:O	3:QC:66:VAL:HA	2.20	0.42
5:QE:74:GLY:O	5:QE:116:THR:N	2.53	0.42
8:QH:12:ARG:NH1	8:QH:27:PRO:HD2	2.35	0.42
14:QN:23:ARG:NH1	14:QN:30:ALA:HB2	2.35	0.42
19:QS:5:LEU:HD21	50:R4:67:TYR:CE2	2.49	0.42
21:QU:2:GLY:O	21:QU:5:ASP:N	2.47	0.42
25:RA:30:G:O2'	25:RA:1214:A:N3	2.45	0.42
25:RA:288:C:H2'	25:RA:289:A:C8	2.54	0.42
25:RA:1338:G:N7	43:RX:62:LYS:NZ	2.67	0.42
25:RA:1430:C:H2'	25:RA:1431:U:C6	2.55	0.42
25:RA:2443:C:H2'	25:RA:2444:G:C8	2.55	0.42
25:RA:2648:C:H2'	25:RA:2649:U:C6	2.55	0.42
27:RD:35:LYS:HB3	27:RD:36:PRO:HA	2.01	0.42
31:RH:76:VAL:C	31:RH:78:GLY:N	2.78	0.42
36:RQ:118:LEU:HD23	36:RQ:118:LEU:HA	1.87	0.42
38:RS:108:GLY:O	38:RS:110:LEU:HG	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:RU:58:ARG:NH1	40:RU:93:LYS:HE2	2.35	0.42
44:RY:39:VAL:HB	44:RY:40:GLU:H	1.57	0.42
49:R3:35:ARG:HB3	49:R3:37:LEU:HD21	2.01	0.42
50:R4:38:LYS:HG3	50:R4:44:THR:OG1	2.20	0.42
52:R6:25:LYS:HD2	54:R8:34:TRP:HZ2	1.84	0.42
1:XA:412:A:H4'	1:XA:413:G:O5'	2.20	0.42
1:XA:1065:U:O2'	1:XA:1066:C:OP2	2.36	0.42
1:XA:1399:C:C4	1:XA:1401:G:C2	3.08	0.42
9:XI:4:TYR:CE1	9:XI:88:TYR:HB2	2.55	0.42
12:XL:89:ARG:HB3	12:XL:97:ARG:HA	2.02	0.42
12:XL:119:LYS:HB2	12:XL:120:TYR:HD1	1.83	0.42
19:XS:40:ILE:CG1	19:XS:41:VAL:HG13	2.47	0.42
25:YA:7:G:H2'	25:YA:8:A:O4'	2.20	0.42
25:YA:1532:C:H2'	25:YA:1533:C:O4'	2.19	0.42
25:YA:1658:C:H2'	25:YA:1659:U:C6	2.55	0.42
25:YA:2277:G:H5"	36:YQ:85:LYS:HB2	2.00	0.42
25:YA:2331:G:H4'	46:Y0:43:THR:H	1.83	0.42
25:YA:2712:U:OP1	25:YA:2714:G:H4'	2.20	0.42
26:YB:15:A:H1'	26:YB:109:G:N9	2.35	0.42
27:YD:145:VAL:O	27:YD:154:LYS:N	2.48	0.42
28:YE:36:ARG:HB3	28:YE:36:ARG:NH1	2.31	0.42
29:YF:20:LEU:HD12	29:YF:21:ALA:N	2.26	0.42
31:YH:105:LEU:CD1	31:YH:105:LEU:N	2.81	0.42
36:YQ:34:LEU:HD23	36:YQ:104:PHE:CD2	2.55	0.42
38:YS:51:ALA:HB3	38:YS:73:LEU:HD23	2.01	0.42
50:Y4:39:CYS:O	50:Y4:40:HIS:HB2	2.20	0.42
1:QA:922:G:C6	1:QA:923:A:C6	3.08	0.41
1:QA:1123:A:O2'	10:QJ:38:ILE:HG22	2.20	0.41
1:QA:1313:U:OP1	19:QS:5:LEU:HB2	2.20	0.41
2:QB:88:ALA:HB2	2:QB:219:VAL:HG13	2.02	0.41
3:QC:134:ILE:HD11	3:QC:153:VAL:HG21	2.02	0.41
3:QC:162:GLN:HE21	3:QC:162:GLN:CA	2.27	0.41
7:QG:13:GLN:O	7:QG:24:THR:HG21	2.20	0.41
8:QH:54:ASP:OD1	8:QH:54:ASP:N	2.53	0.41
8:QH:59:LEU:O	8:QH:61:VAL:HG23	2.20	0.41
25:RA:249:C:O2	54:R8:12:LYS:HE3	2.20	0.41
25:RA:592:G:O2'	54:R8:4:MET:N	2.50	0.41
25:RA:594:U:C5'	54:R8:61:LEU:HD21	2.50	0.41
25:RA:989:G:OP2	49:R3:11:SER:OG	2.25	0.41
27:RD:43:ARG:HB2	27:RD:54:ARG:HB2	2.02	0.41
30:RG:27:ASN:HB3	30:RG:30:GLU:HG3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:RG:47:LYS:HD3	30:RG:81:LYS:CB	2.49	0.41
32:RI:101:LEU:HD13	32:RI:109:ILE:HD13	2.02	0.41
35:RP:98:GLU:HA	35:RP:101:VAL:HG12	2.01	0.41
35:RP:127:ALA:C	35:RP:147:LEU:HD23	2.45	0.41
36:RQ:22:LYS:HA	45:RZ:78:LYS:HD2	2.02	0.41
38:RS:78:LEU:HD23	38:RS:78:LEU:HA	1.86	0.41
45:RZ:109:ALA:C	45:RZ:111:VAL:HG12	2.45	0.41
49:R3:7:LYS:HA	49:R3:33:GLN:O	2.20	0.41
51:R5:40:LYS:HE2	51:R5:47:PRO:CG	2.49	0.41
53:R7:47:ARG:HB2	53:R7:48:LYS:H	1.68	0.41
1:XA:9:G:H5"	5:XE:126:ARG:HE	1.85	0.41
1:XA:405:U:O4	4:XD:2:GLY:N	2.53	0.41
1:XA:502:G:H2'	1:XA:503:C:O4'	2.20	0.41
1:XA:659:U:H2'	1:XA:660:G:C8	2.54	0.41
1:XA:692:U:H1'	1:XA:695:A:N7	2.35	0.41
1:XA:803:G:C6	1:XA:804:U:C4	3.08	0.41
1:XA:994:A:N7	1:XA:1216:G:H4'	2.35	0.41
1:XA:1410:G:H2'	1:XA:1411:C:C6	2.55	0.41
1:XA:1442:G:C5	1:XA:1446:A:C6	3.08	0.41
8:XH:104:ARG:HD2	8:XH:138:TRP:CG	2.55	0.41
25:YA:288:C:H2'	25:YA:289:A:H8	1.84	0.41
25:YA:1348:G:C2'	25:YA:1349:A:H5"	2.48	0.41
25:YA:1810:A:H8	25:YA:1810:A:O5'	2.03	0.41
25:YA:2098:U:H3	25:YA:2191:G:H1	1.68	0.41
25:YA:2393:A:H4'	35:YP:61:ARG:O	2.20	0.41
25:YA:2762:G:C6	25:YA:2763:G:C4	3.08	0.41
27:YD:165:ILE:O	27:YD:166:GLN:NE2	2.53	0.41
28:YE:9:VAL:HB	28:YE:10:GLY:H	1.70	0.41
28:YE:16:ARG:C	28:YE:17:ASP:CG	2.86	0.41
29:YF:53:THR:O	29:YF:55:GLY:N	2.53	0.41
29:YF:101:LEU:HD12	29:YF:102:PRO:N	2.33	0.41
31:YH:84:SER:O	31:YH:85:LYS:CB	2.64	0.41
33:YN:35:ARG:HB2	33:YN:42:TRP:CZ3	2.55	0.41
36:YQ:118:LEU:HD13	36:YQ:131:ILE:HG23	2.02	0.41
37:YR:2:ARG:HG2	37:YR:5:LYS:NZ	2.35	0.41
38:YS:26:LEU:HB3	38:YS:87:PHE:HA	2.02	0.41
43:YX:8:ILE:O	48:Y2:36:ARG:NH2	2.53	0.41
43:YX:26:TYR:HB3	43:YX:92:LEU:HD12	2.02	0.41
44:YY:51:VAL:HG23	44:YY:57:GLN:N	2.36	0.41
47:Y1:76:ARG:H	47:Y1:76:ARG:HD2	1.84	0.41
49:Y3:35:ARG:HB3	49:Y3:37:LEU:HD21	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:8:A:H4'	1:QA:9:G:OP1	2.20	0.41
1:QA:304:U:O2'	1:QA:305:G:H5'	2.20	0.41
1:QA:811:C:H4'	1:QA:900:A:N6	2.34	0.41
1:QA:943:U:H2'	1:QA:944:G:H5'	2.02	0.41
1:QA:1035:A:N6	1:QA:1036:G:N3	2.68	0.41
1:QA:1277:C:O2'	1:QA:1279:A:H8	2.01	0.41
2:QB:208:ILE:HA	2:QB:211:ILE:HD12	2.02	0.41
4:QD:150:GLU:H	4:QD:150:GLU:CD	2.28	0.41
5:QE:79:GLU:HB3	5:QE:92:LYS:HA	2.02	0.41
8:QH:38:ILE:HD12	8:QH:118:VAL:HG12	2.02	0.41
14:QN:29:ARG:HG2	14:QN:31:ARG:O	2.20	0.41
18:QR:56:THR:HB	18:QR:58:LEU:CD1	2.50	0.41
25:RA:152:G:H1	25:RA:174:C:H42	1.67	0.41
25:RA:273:G:H1	25:RA:364:C:H42	1.67	0.41
25:RA:479:A:N3	25:RA:481:G:H5"	2.35	0.41
25:RA:1803:A:H4'	27:RD:259:THR:HG23	2.01	0.41
25:RA:2025:C:H2'	25:RA:2026:C:C6	2.55	0.41
27:RD:101:GLU:OE1	27:RD:103:ARG:NH1	2.53	0.41
31:RH:128:PRO:CG	31:RH:129:THR:H	2.33	0.41
32:RI:33:ARG:HB3	32:RI:35:LEU:HD23	2.02	0.41
33:RN:10:GLU:HA	33:RN:11:PRO:HD3	1.65	0.41
36:RQ:20:ALA:HA	36:RQ:98:LYS:HB3	2.02	0.41
37:RR:109:ALA:HA	37:RR:110:PRO:HD2	1.95	0.41
45:RZ:97:GLU:HB3	45:RZ:125:LEU:HD11	2.02	0.41
47:R1:85:LEU:HA	47:R1:87:PRO:HD2	2.01	0.41
1:XA:390:C:H2'	1:XA:391:G:C8	2.56	0.41
1:XA:1004:A:OP1	1:XA:1025:U:N3	2.53	0.41
1:XA:1024:G:OP1	1:XA:1024:G:H4'	2.21	0.41
1:XA:1194:U:H2'	1:XA:1195:C:C6	2.55	0.41
1:XA:1222:G:P	19:XS:77:THR:HG21	2.60	0.41
1:XA:1336:C:O2'	1:XA:1337:G:P	2.78	0.41
2:XB:24:TRP:CD1	2:XB:24:TRP:H	2.38	0.41
4:XD:127:THR:HA	4:XD:132:ARG:HA	2.03	0.41
9:XI:40:LEU:O	9:XI:42:ARG:N	2.48	0.41
9:XI:46:ALA:HB2	9:XI:74:ILE:HG23	2.01	0.41
10:XJ:3:LYS:HB2	10:XJ:75:ILE:O	2.19	0.41
12:XL:61:THR:C	12:XL:63:GLY:H	2.27	0.41
12:XL:109:GLY:HA3	12:XL:121:GLY:O	2.20	0.41
16:XP:56:ALA:HB1	16:XP:74:LEU:HD13	2.02	0.41
25:YA:155:C:H5'	25:YA:161:U:OP2	2.20	0.41
25:YA:592:G:N3	54:Y8:4:MET:HE2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:658:C:O2'	29:YF:102:PRO:HG3	2.20	0.41
25:YA:1190:G:C5'	35:YP:32:THR:HA	2.49	0.41
25:YA:2210:G:H5'	25:YA:2211:G:C5	2.54	0.41
25:YA:2443:C:OP1	29:YF:68:LYS:HD3	2.20	0.41
25:YA:2712:U:O2'	25:YA:2712(A):A:OP1	2.36	0.41
25:YA:2872:G:O2'	25:YA:2873:A:H5'	2.19	0.41
27:YD:14:ARG:CG	27:YD:15:PHE:N	2.83	0.41
27:YD:109:ASP:HB2	27:YD:197:GLY:HA2	2.03	0.41
28:YE:4:ILE:HG22	28:YE:198:VAL:HB	2.02	0.41
28:YE:167:VAL:CG1	28:YE:189:PRO:HD3	2.50	0.41
28:YE:186:GLY:C	28:YE:188:VAL:H	2.28	0.41
29:YF:64:ILE:HG23	29:YF:65:TRP:CD1	2.54	0.41
30:YG:146:TYR:C	30:YG:148:MET:H	2.28	0.41
32:YI:68:LEU:HA	32:YI:71:ILE:HG22	2.02	0.41
33:YN:59:LYS:HE3	33:YN:61:ARG:HH22	1.85	0.41
37:YR:3:HIS:O	37:YR:5:LYS:N	2.53	0.41
44:YY:84:ARG:O	44:YY:95:LYS:HD3	2.20	0.41
50:Y4:15:ILE:C	50:Y4:16:CYS:SG	3.03	0.41
1:QA:255:G:H1'	17:QQ:16:GLN:NE2	2.35	0.41
1:QA:568:G:O6	12:QL:5:PRO:HD3	2.20	0.41
1:QA:753:A:H4'	1:QA:754:C:O5'	2.20	0.41
3:QC:189:ALA:O	3:QC:190:ARG:C	2.62	0.41
4:QD:146:ILE:H	4:QD:146:ILE:HD12	1.85	0.41
4:QD:166:LYS:HD3	27:YD:134:ARG:NH1	2.35	0.41
19:QS:28:LYS:HA	19:QS:47:HIS:HE1	1.86	0.41
19:QS:50:ALA:HB1	19:QS:57:HIS:HB3	2.02	0.41
25:RA:103:A:O5'	25:RA:103:A:H8	2.03	0.41
25:RA:534:U:H2'	25:RA:535:C:C6	2.55	0.41
25:RA:782:A:O2'	27:RD:225:ALA:HB1	2.21	0.41
25:RA:2047:U:H2'	25:RA:2048:G:H8	1.86	0.41
25:RA:2712:U:O2'	25:RA:2712(A):A:OP1	2.38	0.41
25:RA:2746:U:H5"	31:RH:138:LYS:HE2	2.01	0.41
30:RG:5:VAL:N	50:R4:25:TYR:CE2	2.88	0.41
31:RH:146:ALA:HA	31:RH:164:TYR:OH	2.21	0.41
34:RO:7:TYR:CE1	34:RO:20:MET:HB2	2.56	0.41
34:RO:73:ASP:OD2	39:RT:32:TYR:OH	2.28	0.41
35:RP:121:LYS:HE2	35:RP:121:LYS:HB2	1.75	0.41
41:RV:76:LYS:HB2	41:RV:81:TYR:HB3	2.01	0.41
44:RY:88:LYS:NZ	44:RY:88:LYS:HA	2.35	0.41
50:R4:45:GLY:O	50:R4:46:GLN:C	2.62	0.41
1:XA:1239:A:H62	1:XA:1299:A:H62	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:XE:143:ARG:HE	8:XH:77:GLU:CD	2.29	0.41
12:XL:8:ASN:O	12:XL:11:VAL:HG23	2.20	0.41
16:XP:23:ASP:O	16:XP:26:ARG:HB2	2.20	0.41
20:XT:44:ALA:CB	20:XT:91:LEU:HB2	2.50	0.41
25:YA:2712:U:O2'	25:YA:2712(A):A:P	2.77	0.41
27:YD:145:VAL:CG1	27:YD:146:GLU:N	2.84	0.41
28:YE:35:GLN:HG3	28:YE:37:ARG:NH2	2.35	0.41
29:YF:80:ALA:O	29:YF:83:PHE:HB2	2.20	0.41
31:YH:169:VAL:HG22	31:YH:170:ARG:N	2.26	0.41
33:YN:18:ALA:HB3	33:YN:55:VAL:O	2.19	0.41
35:YP:135:LEU:HD23	35:YP:135:LEU:HA	1.77	0.41
38:YS:83:LYS:HE3	38:YS:84:GLN:CG	2.49	0.41
38:YS:92:TYR:HB2	38:YS:98:VAL:HG11	2.02	0.41
41:YV:22:VAL:HG12	41:YV:23:GLU:H	1.85	0.41
41:YV:72:VAL:HG13	41:YV:85:LYS:HG2	2.01	0.41
46:Y0:36:ILE:HD11	46:Y0:39:ARG:HG2	2.02	0.41
50:Y4:24:THR:OG1	50:Y4:25:TYR:N	2.53	0.41
54:Y8:25:MET:HB3	54:Y8:26:LYS:H	1.69	0.41
1:QA:280:C:C2	17:QQ:38:ARG:HG3	2.56	0.41
1:QA:481:G:O2'	1:QA:482:A:P	2.78	0.41
1:QA:574:A:N3	1:QA:883:C:H1'	2.36	0.41
2:QB:219:VAL:O	2:QB:223:ILE:HG13	2.20	0.41
3:QC:42:LEU:HD12	3:QC:42:LEU:HA	1.79	0.41
5:QE:110:LEU:HD13	5:QE:118:ILE:HG12	2.01	0.41
8:QH:20:TYR:CE2	8:QH:75:ARG:HD2	2.55	0.41
8:QH:36:LEU:HD12	8:QH:59:LEU:HD13	2.02	0.41
12:QL:90:VAL:HG12	12:QL:92:ASP:H	1.85	0.41
12:QL:117:ARG:HB3	12:QL:122:THR:HB	2.02	0.41
13:QM:40:ASN:HA	13:QM:41:PRO:HD3	1.89	0.41
14:QN:4:LYS:O	14:QN:7:ILE:HG12	2.20	0.41
15:QO:31:LEU:O	15:QO:35:ARG:HG3	2.20	0.41
16:QP:20:VAL:HG21	16:QP:32:TYR:CD2	2.56	0.41
16:QP:53:VAL:O	16:QP:57:ARG:HG2	2.21	0.41
19:QS:42:PRO:CG	50:R4:63:TYR:CE2	3.02	0.41
25:RA:270(U):C:H2'	25:RA:270(V):G:H8	1.85	0.41
25:RA:547:A:H3'	25:RA:548:A:C8	2.56	0.41
25:RA:1025:G:C5	25:RA:1135:C:H1'	2.56	0.41
25:RA:1296:G:OP1	25:RA:2709:G:O2'	2.26	0.41
25:RA:2637:U:C4	25:RA:2638:G:C6	3.08	0.41
26:RB:24:G:H5"	26:RB:25:A:OP1	2.20	0.41
28:RE:13:ARG:HH11	28:RE:13:ARG:HB2	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:RE:111:ARG:NE	28:RE:160:TYR:CE1	2.76	0.41
31:RH:105:LEU:CD1	31:RH:105:LEU:N	2.81	0.41
32:RI:97:ILE:HG12	32:RI:97:ILE:H	1.62	0.41
37:RR:33:ARG:HH21	51:R5:55:ARG:HB3	1.85	0.41
46:R0:23:VAL:HA	46:R0:38:VAL:HG22	2.01	0.41
53:R7:1:MET:SD	53:R7:3:ARG:NH2	2.93	0.41
1:XA:25:C:H2'	1:XA:26:A:C8	2.55	0.41
1:XA:59:A:H3'	1:XA:331:G:H22	1.85	0.41
1:XA:411:A:C6	1:XA:429:U:C4	3.09	0.41
1:XA:696:A:H8	1:XA:696:A:O5'	2.04	0.41
1:XA:1073:U:H2'	1:XA:1074:G:C8	2.54	0.41
1:XA:1114:C:H1'	14:XN:60:SER:HB2	2.02	0.41
5:XE:131:ILE:HD13	5:XE:131:ILE:HA	1.84	0.41
9:XI:125:TYR:HD1	9:XI:126:SER:N	2.18	0.41
11:XK:18:ARG:HA	11:XK:81:ASP:H	1.86	0.41
12:XL:62:SER:HB2	12:XL:64:TYR:CD1	2.56	0.41
16:XP:17:TYR:HE2	16:XP:41:PRO:HG3	1.85	0.41
20:XT:11:SER:HA	20:XT:13:LEU:HD12	2.01	0.41
25:YA:137(A):G:H1'	43:YX:41:ASN:ND2	2.36	0.41
25:YA:270:A:OP2	25:YA:270(Y):G:N1	2.41	0.41
25:YA:729:G:OP2	27:YD:13:ARG:NH1	2.53	0.41
25:YA:1059:G:H3'	25:YA:1060:U:H5"	2.02	0.41
25:YA:2146:C:H4'	25:YA:2147:G:C8	2.56	0.41
25:YA:2599:G:OP2	27:YD:236:GLY:HA3	2.19	0.41
25:YA:2764:A:N6	25:YA:2766:G:C2	2.89	0.41
26:YB:66:A:O2'	26:YB:67:G:P	2.78	0.41
27:YD:9:TYR:CZ	27:YD:13:ARG:HD3	2.54	0.41
27:YD:158:ALA:O	27:YD:196:VAL:HG11	2.21	0.41
27:YD:269:PHE:N	27:YD:269:PHE:CD1	2.88	0.41
29:YF:160:ASN:CG	29:YF:163:VAL:HG23	2.45	0.41
29:YF:176:LEU:HD11	29:YF:180:GLY:O	2.19	0.41
30:YG:86:MET:HA	30:YG:87:PRO:HD2	1.95	0.41
30:YG:165:THR:OG1	30:YG:168:GLU:HG3	2.21	0.41
44:YY:63:LYS:HD2	44:YY:63:LYS:HA	1.86	0.41
47:Y1:89:GLU:HA	47:Y1:93:GLU:HB2	2.02	0.41
1:QA:34:C:H2'	1:QA:35:G:H8	1.86	0.41
1:QA:909:A:H2'	1:QA:910:C:O4'	2.20	0.41
2:QB:184:VAL:N	2:QB:198:ASP:OD2	2.44	0.41
11:QK:120:ARG:HA	11:QK:121:PRO:HD3	1.87	0.41
13:QM:65:LYS:HB3	50:R4:50:VAL:HG21	2.03	0.41
13:QM:91:ARG:HB2	13:QM:98:VAL:HG13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:QQ:83:ASP:O	17:QQ:87:LYS:HG2	2.20	0.41
20:QT:89:ARG:HH21	20:QT:104:LEU:HG	1.85	0.41
22:QV:21:A:N6	22:QV:46:G:H2'	2.34	0.41
25:RA:741:G:H2'	25:RA:742:G:H8	1.85	0.41
25:RA:828:U:H4'	25:RA:831:G:N1	2.36	0.41
25:RA:2052:G:H4'	28:RE:143:ASN:O	2.20	0.41
25:RA:2346:A:H5"	25:RA:2383:G:H1'	2.02	0.41
25:RA:2770:G:H5"	25:RA:2771:C:OP2	2.20	0.41
25:RA:2867:G:O2'	25:RA:2868:A:O5'	2.34	0.41
26:RB:27:C:H5'	26:RB:28:C:OP2	2.20	0.41
28:RE:35:GLN:HG3	28:RE:37:ARG:NH2	2.36	0.41
29:RF:135:LYS:HA	29:RF:135:LYS:HD2	1.75	0.41
30:RG:6:ALA:HB3	30:RG:104:GLU:OE2	2.20	0.41
31:RH:86:GLU:HG3	31:RH:165:ALA:CA	2.49	0.41
33:RN:47:ALA:HB2	33:RN:112:LEU:HD11	2.02	0.41
33:RN:71:ILE:HG21	33:RN:84:LYS:HB3	2.02	0.41
35:RP:21:ARG:HB3	35:RP:22:GLY:H	1.61	0.41
36:RQ:34:LEU:HD23	36:RQ:104:PHE:CD2	2.55	0.41
37:RR:10:LEU:O	37:RR:12:ARG:HG3	2.21	0.41
37:RR:22:ARG:HA	37:RR:47:PHE:HE2	1.86	0.41
38:RS:14:VAL:HG21	38:RS:89:ARG:HG2	2.02	0.41
54:R8:26:LYS:HD3	54:R8:26:LYS:HA	1.71	0.41
3:XC:150:LYS:HE3	3:XC:167:TRP:HE1	1.84	0.41
10:XJ:3:LYS:HD2	10:XJ:77:PRO:HD3	2.01	0.41
12:XL:53:ARG:HH12	12:XL:92:ASP:HB3	1.85	0.41
13:XM:12:ASN:N	13:XM:45:VAL:HG13	2.36	0.41
18:XR:38:GLU:O	18:XR:42:ARG:NH1	2.54	0.41
25:YA:83:G:N1	25:YA:102:G:H1'	2.36	0.41
25:YA:376:C:H2'	25:YA:377:C:C6	2.56	0.41
25:YA:588:U:H2'	25:YA:589:C:C6	2.55	0.41
25:YA:892:G:N2	25:YA:893:C:C2	2.88	0.41
25:YA:2017:U:O2	51:Y5:10:LYS:HB2	2.20	0.41
25:YA:2322:A:H2'	25:YA:2323:G:O4'	2.20	0.41
25:YA:2832:U:H4'	25:YA:2833:G:C5'	2.50	0.41
27:YD:168:ARG:O	27:YD:169:GLU:HB2	2.19	0.41
27:YD:182:LEU:N	27:YD:272:ALA:HB3	2.32	0.41
29:YF:129:PHE:O	29:YF:142:TRP:HD1	2.03	0.41
30:YG:166:ASP:OD1	30:YG:166:ASP:N	2.54	0.41
31:YH:146:ALA:HB2	31:YH:164:TYR:OH	2.21	0.41
33:YN:29:LYS:H	33:YN:29:LYS:HG2	1.53	0.41
40:YU:69:CYS:HB3	40:YU:106:PHE:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:YV:38:LEU:O	41:YV:51:VAL:HA	2.20	0.41
46:Y0:72:ARG:HB2	46:Y0:75:LEU:HB2	2.03	0.41
53:Y7:47:ARG:HB2	53:Y7:47:ARG:HE	1.58	0.41
1:QA:771:G:H2'	1:QA:772:U:C6	2.56	0.41
1:QA:793:U:H3'	1:QA:794:A:H5"	2.01	0.41
1:QA:1148:U:OP1	9:QI:7:THR:HG21	2.20	0.41
1:QA:1216:G:H5"	14:QN:5:ALA:HB2	2.03	0.41
2:QB:210:SER:O	2:QB:214:ILE:HG12	2.21	0.41
5:QE:100:VAL:O	5:QE:107:ARG:NH2	2.53	0.41
10:QJ:79:ARG:HD3	10:QJ:79:ARG:HA	1.78	0.41
25:RA:34:C:N4	25:RA:447:A:H61	2.18	0.41
25:RA:244:A:H4'	35:RP:74:GLU:HB2	2.03	0.41
25:RA:372:G:HO2'	25:RA:373:U:H5	1.64	0.41
25:RA:558:G:OP1	33:RN:111:PRO:HD2	2.21	0.41
25:RA:580:C:H2'	25:RA:581:C:C6	2.55	0.41
25:RA:646:A:H2'	25:RA:647:G:O4'	2.19	0.41
25:RA:860:U:H5	25:RA:917:A:C2	2.39	0.41
25:RA:1246:A:OP1	29:RF:38:ARG:NH1	2.50	0.41
25:RA:2115:G:H8	25:RA:2171:A:H62	1.68	0.41
26:RB:13:A:O2'	26:RB:14:U:H3'	2.21	0.41
27:RD:221:VAL:HG22	27:RD:226:MET:HE3	2.03	0.41
28:RE:93:VAL:HG21	28:RE:180:ASN:HA	2.03	0.41
28:RE:167:VAL:CG1	28:RE:189:PRO:HD3	2.50	0.41
28:RE:179:GLU:CB	28:RE:181:LEU:HD23	2.24	0.41
28:RE:197:ILE:HD11	28:RE:199:ARG:NH1	2.30	0.41
30:RG:124:SER:HB2	30:RG:131:TYR:CE1	2.56	0.41
43:RX:51:VAL:HG13	43:RX:81:VAL:HG23	2.03	0.41
44:RY:54:LYS:HB3	44:RY:55:TYR:CE1	2.55	0.41
44:RY:84:ARG:HD3	44:RY:86:ARG:NH1	2.35	0.41
52:R6:13:CYS:O	52:R6:21:TYR:HA	2.20	0.41
54:R8:14:VAL:CG1	54:R8:60:LEU:HD11	2.50	0.41
54:R8:16:ILE:HD11	54:R8:57:ARG:CG	2.44	0.41
1:XA:266:G:O2'	1:XA:267:C:P	2.79	0.41
1:XA:1129:C:H5'	1:XA:1130:A:OP1	2.20	0.41
1:XA:1305:G:H5"	21:XU:4:GLY:C	2.46	0.41
2:XB:80:ILE:HD11	2:XB:208:ILE:HG12	2.02	0.41
15:XO:56:LEU:HD21	25:YA:715:G:C4	2.55	0.41
16:XP:60:LEU:HD23	16:XP:60:LEU:HA	1.81	0.41
17:XQ:59:ILE:HB	17:XQ:71:PHE:HB3	2.03	0.41
20:XT:83:ARG:HA	20:XT:86:ARG:HB3	2.02	0.41
25:YA:704:G:H1'	25:YA:727:A:N6	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:YA:996:A:OP2	40:YU:92:ARG:NH2	2.53	0.41
25:YA:1063:G:H1'	25:YA:1077:A:N7	2.36	0.41
25:YA:1069:A:H4'	25:YA:1070:A:H5"	2.02	0.41
25:YA:1164:G:H1	25:YA:1185:C:H42	1.68	0.41
25:YA:2106:G:H2'	25:YA:2107:C:O4'	2.20	0.41
25:YA:2635:C:H5'	28:YE:77:ILE:CD1	2.51	0.41
25:YA:2726:U:O2'	25:YA:2727:G:H8	2.03	0.41
28:YE:161:GLY:O	28:YE:162:ALA:HB3	2.20	0.41
29:YF:133:ASN:O	29:YF:134:GLY:C	2.64	0.41
30:YG:103:LEU:O	30:YG:107:LEU:HG	2.21	0.41
32:YI:88:ILE:HG12	32:YI:122:GLU:N	2.36	0.41
41:YV:72:VAL:CG1	41:YV:85:LYS:HG2	2.50	0.41
44:YY:32:PRO:C	44:YY:34:LYS:H	2.29	0.41
45:YZ:6:LYS:NZ	45:YZ:43:GLU:HG3	2.36	0.41
50:Y4:48:ARG:NH1	50:Y4:52:THR:H	2.19	0.41
54:Y8:14:VAL:CG1	54:Y8:60:LEU:HD11	2.50	0.41
1:QA:162:A:H8	1:QA:162:A:O5'	2.04	0.41
1:QA:604:G:H2'	1:QA:605:U:O4'	2.20	0.41
1:QA:1280:A:HO2'	1:QA:1281:U:P	2.42	0.41
2:QB:47:THR:HA	2:QB:202:PRO:HG2	2.00	0.41
5:QE:12:LEU:O	5:QE:13:ILE:HD12	2.21	0.41
5:QE:127:ASN:HA	5:QE:128:PRO:HD3	1.89	0.41
12:QL:21:LYS:N	12:QL:21:LYS:CD	2.83	0.41
12:QL:109:GLY:HA3	12:QL:121:GLY:O	2.20	0.41
19:QS:36:ARG:HA	19:QS:71:LEU:HB2	2.02	0.41
20:QT:53:LEU:HD23	20:QT:53:LEU:HA	1.88	0.41
25:RA:220:G:H22	25:RA:427:U:H2'	1.86	0.41
25:RA:445:C:H5"	40:RU:3:ARG:HB3	2.01	0.41
25:RA:508:G:O2'	25:RA:509:C:P	2.78	0.41
25:RA:588:U:H2'	25:RA:589:C:C6	2.55	0.41
25:RA:784:A:O4'	27:RD:227:ASN:ND2	2.54	0.41
25:RA:830:G:N2	25:RA:2445:G:O2'	2.49	0.41
25:RA:2105:C:H2'	25:RA:2106:G:H8	1.85	0.41
25:RA:2420:C:OP1	54:R8:34:TRP:HA	2.21	0.41
26:RB:52:A:O2'	26:RB:53:A:N7	2.50	0.41
28:RE:186:GLY:C	28:RE:188:VAL:H	2.28	0.41
31:RH:45:VAL:O	31:RH:45:VAL:CG1	2.68	0.41
31:RH:137:ASP:HB2	31:RH:140:LYS:CE	2.51	0.41
50:R4:68:ARG:HB2	50:R4:69:LYS:H	1.35	0.41
54:R8:64:TYR:HB3	54:R8:65:GLU:H	1.40	0.41
1:XA:8:A:N6	4:XD:208:SER:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:254:G:OP1	17:XQ:67:LYS:O	2.39	0.41
1:XA:632:A:H8	1:XA:633:G:C8	2.37	0.41
1:XA:749:C:H2'	1:XA:750:G:H8	1.85	0.41
1:XA:947:G:H2'	1:XA:948:C:O4'	2.21	0.41
1:XA:1118:C:H1'	1:XA:1179:A:C4	2.56	0.41
1:XA:1173:G:H2'	1:XA:1174:G:O4'	2.21	0.41
1:XA:1226:C:O2'	13:XM:103:THR:O	2.32	0.41
1:XA:1301:U:HO2'	1:XA:1302:U:P	2.42	0.41
1:XA:1390:U:H2'	1:XA:1391:U:C6	2.55	0.41
2:XB:118:LEU:CB	2:XB:142:LEU:HD12	2.50	0.41
8:XH:44:PHE:HE2	8:XH:109:ILE:CG2	2.34	0.41
12:XL:25:PRO:HD2	12:XL:97:ARG:HH11	1.86	0.41
12:XL:90:VAL:HG12	12:XL:92:ASP:H	1.85	0.41
14:XN:51:GLY:O	14:XN:53:LEU:N	2.53	0.41
18:XR:85:LEU:HD23	18:XR:88:LYS:HD2	2.03	0.41
25:YA:242:G:H5'	54:Y8:62:LEU:CD2	2.40	0.41
25:YA:404:C:O2'	25:YA:405:U:OP2	2.31	0.41
25:YA:675:A:OP1	29:YF:63:LYS:NZ	2.53	0.41
25:YA:1085:A:O2'	25:YA:1086:A:P	2.79	0.41
25:YA:1263:U:O2'	51:Y5:11:THR:HG23	2.21	0.41
25:YA:1614:A:N6	42:YW:88:ARG:H	2.18	0.41
25:YA:1799:G:H4'	25:YA:1800:C:O5'	2.21	0.41
25:YA:1819:A:H4'	25:YA:1820:U:O5'	2.21	0.41
25:YA:2680:C:OP1	28:YE:109:LYS:HG3	2.21	0.41
25:YA:2693:A:H2'	25:YA:2694:G:H8	1.86	0.41
29:YF:68:LYS:O	29:YF:69:HIS:HB2	2.20	0.41
31:YH:76:VAL:C	31:YH:78:GLY:N	2.78	0.41
31:YH:86:GLU:HG3	31:YH:165:ALA:CA	2.49	0.41
32:YI:81:VAL:HG21	32:YI:88:ILE:HD12	2.03	0.41
38:YS:66:ALA:HA	38:YS:69:VAL:CG1	2.51	0.41
38:YS:93:LYS:HE3	38:YS:93:LYS:HB2	1.93	0.41
40:YU:98:LEU:HD23	40:YU:99:ALA:N	2.36	0.41
48:Y2:61:LEU:HD23	48:Y2:61:LEU:HA	1.85	0.41
54:Y8:26:LYS:HD3	54:Y8:26:LYS:HA	1.86	0.41
1:QA:37:U:O2'	1:QA:500:G:H4'	2.21	0.41
1:QA:137:C:N4	1:QA:226:G:H1	2.19	0.41
1:QA:145:G:H2'	1:QA:146:G:O4'	2.21	0.41
1:QA:828:A:H2'	1:QA:829:G:O4'	2.21	0.41
1:QA:1077:G:N2	1:QA:1080:A:OP2	2.51	0.41
3:QC:122:GLU:HA	3:QC:125:GLU:OE1	2.21	0.41
4:QD:96:LEU:HD13	4:QD:96:LEU:HA	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:QD:166:LYS:HG3	4:QD:178:VAL:HG11	2.02	0.41
5:QE:143:ARG:HH21	8:QH:77:GLU:CD	2.29	0.41
7:QG:45:ASP:O	7:QG:48:LYS:HB3	2.21	0.41
9:QI:40:LEU:HD11	9:QI:70:LYS:HG2	2.02	0.41
12:QL:43:VAL:HG13	12:QL:55:VAL:HG21	2.03	0.41
12:QL:53:ARG:HH12	12:QL:92:ASP:CB	2.33	0.41
22:QV:2:C:C2'	22:QV:3:G:H5'	2.50	0.41
25:RA:18:C:H2'	25:RA:19:C:C6	2.56	0.41
25:RA:277:C:H3'	25:RA:278:A:H5'	2.01	0.41
25:RA:530:G:C5	25:RA:2022:U:H5"	2.55	0.41
25:RA:634:C:H2'	25:RA:635:C:H6	1.86	0.41
25:RA:686:G:H21	25:RA:788:A:H61	1.69	0.41
25:RA:917:A:O2'	26:RB:98:G:H4'	2.21	0.41
25:RA:1061:U:H3'	25:RA:1062:G:H5"	2.02	0.41
25:RA:2516:G:C6	25:RA:2517:C:C4	3.09	0.41
25:RA:2836:U:H2'	25:RA:2837:G:C8	2.56	0.41
25:RA:2893:G:H5"	25:RA:2894:G:C5'	2.50	0.41
27:RD:127:VAL:HA	27:RD:193:VAL:HG22	2.02	0.41
27:RD:226:MET:HB3	27:RD:230:ASP:HB2	2.03	0.41
28:RE:4:ILE:HG22	28:RE:198:VAL:HB	2.02	0.41
28:RE:71:GLY:O	28:RE:72:VAL:C	2.63	0.41
28:RE:134:ILE:C	28:RE:134:ILE:CD1	2.94	0.41
28:RE:179:GLU:C	28:RE:181:LEU:H	2.28	0.41
30:RG:47:LYS:HB2	30:RG:47:LYS:HE3	1.81	0.41
30:RG:179:PRO:CG	50:R4:38:LYS:HZ2	2.15	0.41
33:RN:57:ALA:O	33:RN:60:ILE:HD11	2.21	0.41
33:RN:73:THR:HB	33:RN:82:LEU:HD11	2.02	0.41
34:RO:26:LYS:HB2	34:RO:30:ALA:HB2	2.02	0.41
36:RQ:20:ALA:HB1	36:RQ:99:PRO:CG	2.51	0.41
36:RQ:139:GLU:CG	36:RQ:140:ALA:N	2.84	0.41
1:XA:427:U:C4	1:XA:428:G:C6	3.09	0.41
1:XA:514:C:H2'	1:XA:515:G:C8	2.56	0.41
1:XA:673:G:H5"	6:XF:87:ARG:NH1	2.36	0.41
1:XA:674:G:H2'	1:XA:675:A:C8	2.56	0.41
1:XA:922:G:H1	1:XA:1395:C:H42	1.67	0.41
1:XA:1046:A:H3'	1:XA:1047:G:H8	1.86	0.41
1:XA:1129:C:H4'	1:XA:1130:A:H8	1.85	0.41
1:XA:1211:U:H4'	1:XA:1213:A:H1'	2.02	0.41
1:XA:1415:G:C6	1:XA:1486:G:C6	3.09	0.41
4:XD:64:LEU:HD13	4:XD:198:VAL:HG11	2.02	0.41
6:XF:46:ARG:HB3	6:XF:60:PHE:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:XL:62:SER:O	12:XL:64:TYR:N	2.54	0.41
17:XQ:62:SER:CB	17:XQ:72:ARG:HE	2.33	0.41
20:XT:74:LYS:C	20:XT:76:ALA:H	2.29	0.41
25:YA:1328:G:H2'	25:YA:1330:C:C5	2.56	0.41
25:YA:1534:G:N2	25:YA:1537:C:O2	2.54	0.41
25:YA:1537:C:H2'	25:YA:1538:G:O4'	2.21	0.41
25:YA:1991:U:H2'	25:YA:1992:G:H5"	2.02	0.41
25:YA:2390:U:O2'	25:YA:2391:G:H5'	2.20	0.41
25:YA:2394:C:OP1	35:YP:63:PRO:HD2	2.20	0.41
25:YA:2466:C:OP1	55:Y9:4:ARG:HB2	2.20	0.41
25:YA:2819:G:H2'	25:YA:2821:A:N7	2.36	0.41
27:YD:117:VAL:HG22	27:YD:118:VAL:N	2.35	0.41
27:YD:197:GLY:O	27:YD:198:ASN:HB3	2.21	0.41
28:YE:24:THR:HB	28:YE:184:VAL:HG23	2.02	0.41
28:YE:63:LEU:CD1	28:YE:64:LYS:N	2.71	0.41
29:YF:28:ILE:HD12	29:YF:28:ILE:C	2.46	0.41
30:YG:101:ILE:HD13	50:Y4:9:LEU:HD11	2.02	0.41
30:YG:124:SER:HB2	30:YG:131:TYR:CE1	2.56	0.41
31:YH:53:GLU:OE1	31:YH:53:GLU:HA	2.16	0.41
32:YI:130:TYR:C	32:YI:131:LYS:HD2	2.45	0.41
33:YN:96:GLU:O	33:YN:100:GLU:HG3	2.20	0.41
35:YP:3:LEU:HD23	35:YP:3:LEU:HA	1.94	0.41
40:YU:8:VAL:O	40:YU:12:ARG:HG3	2.20	0.41
40:YU:17:ILE:HG23	40:YU:39:LEU:HD12	2.02	0.41
45:YZ:77:ASP:OD2	45:YZ:80:ARG:HD3	2.21	0.41
45:YZ:97:GLU:HG3	45:YZ:127:LYS:NZ	2.35	0.41
1:QA:498:A:H4'	1:QA:500:G:OP1	2.21	0.41
1:QA:510:A:OP2	4:QD:49:ARG:NH2	2.53	0.41
1:QA:616:G:H2'	1:QA:617:G:H8	1.86	0.41
1:QA:766:A:H2'	1:QA:767:A:O4'	2.20	0.41
1:QA:807:A:H2'	1:QA:808:C:C6	2.56	0.41
1:QA:869:G:H4'	1:QA:872:A:O4'	2.21	0.41
1:QA:939:G:H5"	7:QG:102:ARG:NH2	2.36	0.41
1:QA:1143:G:H2'	1:QA:1144:G:C8	2.55	0.41
1:QA:1187:G:C6	1:QA:1188:A:C5	3.08	0.41
1:QA:1281:U:OP2	1:QA:1282:C:H5	2.03	0.41
1:QA:1417:G:N2	1:QA:1482:G:H2'	2.36	0.41
3:QC:71:ALA:HB2	3:QC:109:PRO:HB3	2.03	0.41
5:QE:47:LYS:HE2	5:QE:47:LYS:HB2	1.82	0.41
8:QH:41:ARG:NH2	8:QH:123:GLU:OE2	2.54	0.41
11:QK:25:TYR:CZ	11:QK:87:THR:HB	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:QK:124:LYS:C	11:QK:126:ARG:H	2.29	0.41
12:QL:8:ASN:O	12:QL:11:VAL:HG23	2.20	0.41
19:QS:23:ASN:C	19:QS:25:LYS:H	2.29	0.41
25:RA:813:U:H2'	25:RA:814:C:C6	2.56	0.41
25:RA:968:G:H2'	25:RA:969:U:O4'	2.21	0.41
25:RA:996:A:H4'	40:RU:92:ARG:NE	2.29	0.41
25:RA:1065:U:H3	25:RA:1073:A:H61	1.69	0.41
25:RA:1130:U:HO2'	25:RA:1131:G:P	2.44	0.41
25:RA:1142(A):A:H4'	33:RN:25:ARG:HH22	1.86	0.41
25:RA:1291:C:H2'	25:RA:1292:U:C6	2.56	0.41
25:RA:1323:U:OP1	42:RW:98:LYS:NZ	2.33	0.41
25:RA:1356:G:C6	25:RA:1357:U:C4	3.09	0.41
25:RA:1519:G:C6	25:RA:1520:U:C4	3.09	0.41
25:RA:1632:A:C6	25:RA:1633:G:C6	3.09	0.41
25:RA:1639:U:H4'	25:RA:2699:C:H4'	2.02	0.41
25:RA:2052:G:OP1	28:RE:141:ILE:HG12	2.20	0.41
25:RA:2150:U:H2'	25:RA:2151:G:C8	2.50	0.41
25:RA:2350:C:H5	54:R8:42:ARG:CZ	2.34	0.41
25:RA:2445:G:OP1	29:RF:74:ARG:NH2	2.54	0.41
25:RA:2466:C:C2	25:RA:2485:G:C2	3.09	0.41
25:RA:2656:U:O4	25:RA:2665:A:H2	2.02	0.41
25:RA:2693:A:H2'	25:RA:2694:G:H8	1.86	0.41
25:RA:2867:G:O2'	25:RA:2868:A:C8	2.72	0.41
27:RD:111:LEU:HD23	27:RD:111:LEU:HA	1.78	0.41
27:RD:150:LYS:N	27:RD:150:LYS:HD3	2.36	0.41
27:RD:257:LEU:HD23	27:RD:257:LEU:HA	1.90	0.41
28:RE:51:PHE:CG	28:RE:52:LEU:N	2.89	0.41
28:RE:161:GLY:O	28:RE:162:ALA:HB3	2.20	0.41
29:RF:78:ILE:H	29:RF:78:ILE:HG13	1.73	0.41
30:RG:131:TYR:O	30:RG:159:VAL:HG13	2.21	0.41
31:RH:145:ALA:O	31:RH:148:ILE:HB	2.21	0.41
36:RQ:80:GLU:C	36:RQ:81:VAL:CG1	2.89	0.41
36:RQ:139:GLU:HG2	36:RQ:140:ALA:N	2.36	0.41
38:RS:62:LYS:HB3	38:RS:97:ARG:CD	2.44	0.41
39:RT:91:ARG:HB2	39:RT:121:ILE:HG13	2.03	0.41
40:RU:61:TRP:O	40:RU:65:ILE:HG13	2.21	0.41
40:RU:83:LEU:HG	40:RU:88:ILE:HB	2.03	0.41
45:RZ:76:LEU:HD23	45:RZ:76:LEU:H	1.86	0.41
47:R1:58:ILE:CD1	47:R1:86:SER:HB2	2.50	0.41
50:R4:14:ILE:HA	50:R4:31:ILE:O	2.21	0.41
50:R4:48:ARG:O	50:R4:49:PHE:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:R6:24:GLU:HB3	52:R6:25:LYS:H	1.74	0.41
54:R8:3:LYS:HE2	54:R8:3:LYS:HB3	1.82	0.41
1:XA:380:G:N2	1:XA:383:A:OP2	2.53	0.41
1:XA:392:G:H2'	1:XA:393:A:C8	2.56	0.41
1:XA:567:G:H2'	1:XA:568:G:O4'	2.20	0.41
1:XA:584:G:H2'	1:XA:585:G:C8	2.55	0.41
1:XA:660:G:H2'	1:XA:661:G:O4'	2.20	0.41
1:XA:1120:G:H2'	1:XA:1121:U:H6	1.84	0.41
1:XA:1129:C:H4'	1:XA:1130:A:H5'	2.03	0.41
2:XB:48:MET:HE2	2:XB:48:MET:HB3	1.94	0.41
3:XC:119:ARG:HH21	3:XC:140:ARG:CZ	2.34	0.41
3:XC:138:VAL:HG13	3:XC:149:ALA:HB3	2.02	0.41
4:XD:52:SER:O	4:XD:55:ALA:HB3	2.21	0.41
4:XD:131:ARG:H	4:XD:131:ARG:HG2	1.67	0.41
8:XH:6:ILE:HB	8:XH:85:ARG:NH1	2.36	0.41
9:XI:79:LEU:O	9:XI:83:ARG:HG2	2.19	0.41
10:XJ:76:ASN:HA	10:XJ:77:PRO:HD2	1.96	0.41
11:XK:19:ALA:CB	11:XK:32:ILE:HG22	2.50	0.41
11:XK:48:ILE:HD13	11:XK:48:ILE:HA	1.83	0.41
12:XL:22:SER:C	12:XL:24:VAL:N	2.79	0.41
12:XL:38:THR:HG22	12:XL:57:LYS:HB3	2.01	0.41
12:XL:117:ARG:HB3	12:XL:122:THR:HB	2.02	0.41
25:YA:195:A:H5"	25:YA:196:A:O5'	2.21	0.41
25:YA:270(H):C:H2'	25:YA:270(I):G:C8	2.56	0.41
25:YA:385:C:O2	35:YP:71:VAL:HG21	2.20	0.41
25:YA:414:C:H2'	25:YA:415:A:C8	2.55	0.41
25:YA:646:A:H2'	25:YA:647:G:O4'	2.21	0.41
25:YA:738:G:C6	25:YA:739:G:C2	3.09	0.41
25:YA:831:G:O2'	35:YP:38:GLN:OE1	2.34	0.41
25:YA:860:U:C5	25:YA:917:A:C2	3.08	0.41
25:YA:953:A:OP1	36:YQ:18:LYS:HE3	2.20	0.41
25:YA:1162:G:H4'	41:YV:24:LYS:HB3	2.02	0.41
25:YA:1359:A:C8	25:YA:1359:A:C5'	3.04	0.41
25:YA:1520:U:H2'	25:YA:1521:G:O4'	2.21	0.41
25:YA:1588:C:H2'	25:YA:1589:C:C6	2.56	0.41
25:YA:2016:U:H2'	25:YA:2017:U:O4'	2.20	0.41
25:YA:2498:C:O2'	25:YA:2499:C:H5'	2.21	0.41
25:YA:2849:U:OP2	39:YT:95:ARG:NH1	2.54	0.41
25:YA:2853:C:H2'	25:YA:2854:G:H8	1.85	0.41
26:YB:51:G:N7	38:YS:62:LYS:NZ	2.57	0.41
27:YD:147:LEU:CD1	27:YD:155:LEU:HD21	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:YE:134:ILE:C	28:YE:134:ILE:CD1	2.93	0.41
29:YF:13:SER:OG	29:YF:14:PRO:HD2	2.21	0.41
30:YG:61:ALA:O	50:Y4:7:PRO:HG3	2.21	0.41
30:YG:109:VAL:HG21	50:Y4:14:ILE:CD1	2.51	0.41
31:YH:137:ASP:HB2	31:YH:140:LYS:CE	2.51	0.41
34:YO:4:PRO:O	34:YO:5:GLN:CB	2.69	0.41
35:YP:1:MET:HB3	35:YP:2:LYS:H	1.72	0.41
35:YP:100:LEU:HD13	35:YP:100:LEU:HA	1.87	0.41
36:YQ:139:GLU:CD	36:YQ:139:GLU:H	2.28	0.41
37:YR:44:LEU:HD22	37:YR:48:VAL:HG23	2.03	0.41
38:YS:20:ARG:HE	38:YS:21:THR:HA	1.86	0.41
38:YS:53:SER:HA	38:YS:56:LEU:CD2	2.50	0.41
38:YS:61:ASN:OD1	38:YS:63:THR:OG1	2.29	0.41
38:YS:88:ASP:CG	38:YS:89:ARG:H	2.28	0.41
40:YU:96:ALA:HA	40:YU:98:LEU:HD23	2.03	0.41
43:YX:31:HIS:HB3	43:YX:34:ALA:HB2	2.03	0.41
45:YZ:152:ALA:HB2	45:YZ:168:GLU:HA	2.03	0.41
48:Y2:41:ILE:HD12	48:Y2:43:GLN:N	2.35	0.41
54:Y8:17:THR:O	54:Y8:20:GLY:N	2.47	0.41
55:Y9:2:LYS:HA	55:Y9:2:LYS:HD2	1.86	0.41
1:QA:41:G:H2'	1:QA:42:G:C8	2.55	0.41
1:QA:343:U:O2	1:QA:346:G:N2	2.36	0.41
1:QA:347:G:O2'	1:QA:348:G:OP2	2.30	0.41
1:QA:976:G:OP2	1:QA:1358:U:O2'	2.39	0.41
1:QA:1359:C:O2'	1:QA:1361:G:N7	2.53	0.41
1:QA:1392:G:N2	1:QA:1502:A:H8	2.12	0.41
3:QC:83:ARG:O	3:QC:86:VAL:HG22	2.21	0.41
5:QE:147:ASP:O	5:QE:151:LEU:HG	2.21	0.41
6:QF:35:ALA:HA	6:QF:67:MET:HB3	2.02	0.41
8:QH:44:PHE:HD1	8:QH:80:ILE:HG12	1.86	0.41
11:QK:99:GLN:HG2	11:QK:105:VAL:CG2	2.44	0.41
12:QL:62:SER:HB2	12:QL:64:TYR:CD1	2.56	0.41
20:QT:86:ARG:O	20:QT:90:GLN:HG3	2.21	0.41
25:RA:17:G:H2'	25:RA:18:C:C6	2.56	0.41
25:RA:229:A:OP1	25:RA:229:A:H4'	2.20	0.41
25:RA:273(E):U:H2'	25:RA:273(F):C:O4'	2.20	0.41
25:RA:389:G:N1	35:RP:70:GLN:HB3	2.36	0.41
25:RA:594:U:P	54:R8:61:LEU:HD22	2.60	0.41
25:RA:827:U:H1'	25:RA:2246:G:O2'	2.21	0.41
25:RA:921:G:H4'	25:RA:2269:A:C5	2.56	0.41
25:RA:1534:G:H2'	25:RA:1535:U:H4'	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:RA:1657:C:H2'	25:RA:1658:C:C6	2.56	0.41
25:RA:2115:G:H8	25:RA:2171:A:N6	2.19	0.41
25:RA:2135:A:H62	25:RA:2156:G:N2	2.09	0.41
25:RA:2638:G:OP2	28:RE:82:ARG:NH2	2.54	0.41
25:RA:2699:C:H2'	25:RA:2700:C:O4'	2.21	0.41
28:RE:65:GLY:O	28:RE:66:HIS:C	2.64	0.41
28:RE:147:PRO:HB2	28:RE:149:ARG:HG2	2.03	0.41
29:RF:60:SER:OG	29:RF:61:GLY:N	2.54	0.41
29:RF:167:ALA:HB1	29:RF:173:VAL:HG11	2.03	0.41
50:R4:12:ALA:HB1	50:R4:30:GLU:N	2.34	0.41
50:R4:42:PHE:CZ	50:R4:43:TYR:HB3	2.56	0.41
1:XA:1032(A):G:H2'	1:XA:1032(B):G:H8	1.84	0.41
2:XB:126:GLU:O	2:XB:129:GLU:HB2	2.20	0.41
4:XD:186:LEU:HD23	4:XD:186:LEU:HA	1.95	0.41
6:XF:25:ILE:HD13	6:XF:28:ARG:NH1	2.35	0.41
8:XH:12:ARG:HD3	8:XH:26:VAL:HB	2.03	0.41
10:XJ:6:ILE:O	10:XJ:71:LEU:HD12	2.21	0.41
25:YA:676:A:C2	25:YA:677:A:C8	3.08	0.41
25:YA:849:A:N1	49:Y3:25:ALA:HB2	2.36	0.41
25:YA:1022:G:O2'	25:YA:1023:U:OP2	2.34	0.41
25:YA:1047:G:H2'	25:YA:1110:G:C2	2.56	0.41
25:YA:1061:U:H3'	25:YA:1062:G:C5'	2.51	0.41
25:YA:2002:G:H2'	25:YA:2003:G:C8	2.56	0.41
27:YD:68:LYS:HG3	27:YD:68:LYS:O	2.20	0.41
27:YD:107:ALA:HA	27:YD:108:PRO:HD2	2.01	0.41
28:YE:179:GLU:C	28:YE:181:LEU:H	2.28	0.41
30:YG:7:LEU:HD12	30:YG:104:GLU:HA	2.03	0.41
30:YG:105:LYS:NZ	50:Y4:26:SER:HB2	2.36	0.41
31:YH:20:ALA:HB3	31:YH:23:ARG:HG2	2.03	0.41
31:YH:146:ALA:HA	31:YH:164:TYR:OH	2.21	0.41
36:YQ:60:ARG:HA	45:YZ:178:GLU:C	2.46	0.41
36:YQ:139:GLU:CG	36:YQ:140:ALA:N	2.84	0.41
38:YS:89:ARG:HG2	38:YS:89:ARG:NH1	2.36	0.41
44:YY:67:LEU:O	44:YY:68:HIS:C	2.63	0.41
48:Y2:11:GLU:HA	48:Y2:14:ARG:HD2	2.02	0.41
52:Y6:36:LEU:HD13	52:Y6:50:ARG:CZ	2.51	0.41
54:Y8:38:GLY:C	54:Y8:40:GLU:N	2.79	0.41
54:Y8:53:PRO:C	54:Y8:55:ALA:N	2.76	0.41
1:QA:329:A:C5	1:QA:332:G:C6	3.09	0.40
1:QA:652:U:H1'	1:QA:653:A:H2	1.85	0.40
1:QA:868:C:H2'	1:QA:869:G:O4'	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:1285:A:H4'	1:QA:1286:A:O5'	2.21	0.40
1:QA:1314:C:P	19:QS:6:LYS:HD2	2.61	0.40
5:QE:9:LYS:HB3	5:QE:112:LEU:HD11	2.04	0.40
7:QG:54:THR:O	7:QG:56:GLN:N	2.52	0.40
10:QJ:51:ARG:HG3	10:QJ:60:ARG:C	2.45	0.40
10:QJ:61:GLU:OE1	14:QN:58:LYS:HE2	2.21	0.40
12:QL:62:SER:C	12:QL:64:TYR:N	2.79	0.40
12:QL:89:ARG:HB3	12:QL:97:ARG:HA	2.02	0.40
25:RA:744:G:OP1	28:RE:132:HIS:HA	2.21	0.40
25:RA:774:A:HO2'	25:RA:775:G:P	2.43	0.40
25:RA:969:U:H2'	25:RA:970:C:C6	2.56	0.40
25:RA:1007:C:OP1	33:RN:35:ARG:NH1	2.54	0.40
25:RA:1916:A:H2'	25:RA:1917:U:O4'	2.22	0.40
28:RE:70:ALA:O	28:RE:71:GLY:C	2.62	0.40
33:RN:9:VAL:HG21	33:RN:48:MET:HB3	2.02	0.40
34:RO:63:VAL:HB	34:RO:106:LEU:HD11	2.02	0.40
36:RQ:39:PRO:HA	36:RQ:97:VAL:O	2.21	0.40
36:RQ:76:LYS:HB3	36:RQ:90:VAL:CG1	2.51	0.40
47:R1:95:LEU:HD23	47:R1:95:LEU:HA	1.94	0.40
50:R4:63:TYR:O	50:R4:65:ASP:N	2.54	0.40
1:XA:20:U:H2'	1:XA:21:G:O4'	2.21	0.40
1:XA:388:G:O2'	1:XA:389:A:P	2.79	0.40
1:XA:593:G:H2'	1:XA:594:G:O4'	2.22	0.40
3:XC:42:LEU:HD12	3:XC:42:LEU:HA	1.87	0.40
4:XD:22:LYS:HB3	4:XD:22:LYS:HZ3	1.86	0.40
19:XS:36:ARG:HA	19:XS:71:LEU:HB2	2.03	0.40
20:XT:75:ASN:OD1	20:XT:75:ASN:N	2.40	0.40
25:YA:153:C:OP1	47:Y1:88:LYS:HE2	2.21	0.40
25:YA:565:C:H4'	25:YA:1253:A:C6	2.56	0.40
25:YA:1225:C:O2'	41:YV:85:LYS:HA	2.21	0.40
25:YA:2286:A:H2'	52:Y6:31:PRO:HG2	2.02	0.40
25:YA:2298:A:H2'	25:YA:2299:G:O4'	2.21	0.40
27:YD:228:PRO:HD3	27:YD:234:GLY:O	2.21	0.40
29:YF:62:ARG:HB3	29:YF:62:ARG:CZ	2.51	0.40
29:YF:118:ALA:HA	29:YF:123:LEU:HB3	2.02	0.40
29:YF:198:ALA:HA	29:YF:201:VAL:CG1	2.41	0.40
30:YG:18:GLU:OE1	30:YG:22:ARG:NH1	2.49	0.40
31:YH:26:VAL:HG12	31:YH:33:LEU:HB2	2.03	0.40
31:YH:45:VAL:O	31:YH:45:VAL:CG1	2.69	0.40
31:YH:145:ALA:O	31:YH:148:ILE:HB	2.21	0.40
32:YI:40:THR:O	32:YI:44:LEU:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:YI:131:LYS:HB3	32:YI:132:PRO:HA	2.01	0.40
35:YP:115:LEU:HB2	35:YP:116:GLY:H	1.53	0.40
48:Y2:37:PHE:O	48:Y2:40:SER:HB3	2.21	0.40
48:Y2:53:LEU:O	48:Y2:57:ILE:HG13	2.21	0.40
50:Y4:43:TYR:O	50:Y4:46:GLN:HA	2.20	0.40
50:Y4:48:ARG:HH12	50:Y4:52:THR:H	1.68	0.40
54:Y8:40:GLU:O	54:Y8:42:ARG:N	2.54	0.40
1:QA:173:U:H5"	1:QA:197:A:O4'	2.21	0.40
1:QA:542:G:H5"	4:QD:10:ARG:HH22	1.86	0.40
1:QA:1045:C:H2'	1:QA:1046:A:O4'	2.21	0.40
1:QA:1298:C:N4	7:QG:114:ARG:HD2	2.36	0.40
6:QF:33:TYR:CE1	6:QF:78:GLU:HG2	2.56	0.40
7:QG:93:PRO:O	7:QG:96:GLN:HB2	2.21	0.40
10:QJ:4:ILE:HA	10:QJ:100:THR:HG22	2.02	0.40
12:QL:45:PRO:C	12:QL:46:LYS:O	2.63	0.40
13:QM:40:ASN:O	13:QM:43:THR:OG1	2.29	0.40
19:QS:5:LEU:HD12	19:QS:5:LEU:HA	1.93	0.40
21:QU:10:ARG:HA	21:QU:13:ILE:HB	2.02	0.40
25:RA:848:G:O6	25:RA:929:G:H2'	2.20	0.40
25:RA:861:A:H2'	25:RA:862:G:O4'	2.21	0.40
25:RA:1316:U:H2'	25:RA:1317:A:H8	1.87	0.40
25:RA:1512:G:H2'	25:RA:1513:C:C6	2.57	0.40
25:RA:1665:A:H4'	34:RO:67:LYS:HB2	2.03	0.40
25:RA:1728:G:C6	25:RA:1730:U:OP2	2.74	0.40
25:RA:2198:A:C2	32:RI:29:TYR:HB2	2.56	0.40
25:RA:2566:A:H4'	25:RA:2567:G:O5'	2.21	0.40
25:RA:2742:C:OP1	55:R9:35:ARG:HD3	2.20	0.40
27:RD:44:ASN:HB2	27:RD:49:ILE:HA	2.02	0.40
27:RD:96:HIS:CD2	27:RD:102:LYS:HG2	2.56	0.40
30:RG:59:GLU:O	30:RG:63:ILE:HG23	2.20	0.40
31:RH:26:VAL:HG12	31:RH:33:LEU:HB2	2.03	0.40
32:RI:48:GLU:OE1	32:RI:52:ARG:NH2	2.54	0.40
32:RI:52:ARG:O	32:RI:56:LYS:CB	2.69	0.40
32:RI:128:LEU:HD23	32:RI:140:LEU:HD21	2.03	0.40
38:RS:23:ARG:HB2	38:RS:86:ALA:HB2	2.03	0.40
39:RT:51:ARG:CG	39:RT:98:LYS:HG3	2.52	0.40
39:RT:109:GLU:O	39:RT:113:LYS:HB2	2.21	0.40
46:R0:12:ASN:HB3	46:R0:13:GLY:H	1.67	0.40
46:R0:72:ARG:HB2	46:R0:75:LEU:HB2	2.03	0.40
1:XA:411:A:C4	1:XA:413:G:H1'	2.56	0.40
1:XA:520:A:O2'	12:XL:73:GLU:HG2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:XA:601:C:H2'	1:XA:602:A:H8	1.87	0.40
1:XA:679:C:H2'	1:XA:680:C:C6	2.57	0.40
3:XC:20:SER:HB2	3:XC:40:ARG:NH2	2.29	0.40
3:XC:188:LEU:HD13	3:XC:188:LEU:HA	1.90	0.40
4:XD:165:MET:O	4:XD:167:GLY:N	2.54	0.40
5:XE:12:LEU:HB3	5:XE:31:LEU:CB	2.51	0.40
5:XE:51:VAL:HB	5:XE:52:PRO:HD3	2.04	0.40
9:XI:95:LYS:NZ	9:XI:96:LEU:HD13	2.36	0.40
10:XJ:61:GLU:HG3	14:XN:58:LYS:HE2	2.03	0.40
25:YA:631:A:OP1	35:YP:64:LYS:HE2	2.21	0.40
25:YA:1733:G:H5'	25:YA:1734:C:OP2	2.22	0.40
25:YA:1849:G:H2'	25:YA:1850:G:H8	1.86	0.40
25:YA:2031:A:N6	25:YA:2498:C:O4'	2.55	0.40
25:YA:2853:C:H2'	25:YA:2854:G:C8	2.57	0.40
25:YA:2875:C:O2'	39:YT:5:ALA:HB3	2.21	0.40
27:YD:31:LYS:O	27:YD:32:SER:O	2.39	0.40
27:YD:134:ARG:H	27:YD:134:ARG:HG3	1.55	0.40
28:YE:71:GLY:O	28:YE:72:VAL:C	2.63	0.40
31:YH:65:HIS:O	31:YH:66:GLY:C	2.63	0.40
31:YH:170:ARG:HB3	31:YH:171:LEU:H	1.47	0.40
52:Y6:11:LEU:HD13	52:Y6:11:LEU:HA	1.86	0.40
52:Y6:13:CYS:HB2	52:Y6:22:ALA:HB3	2.03	0.40
1:QA:9:G:H2'	1:QA:10:A:H8	1.86	0.40
1:QA:883:C:O2'	1:QA:884:U:H5'	2.22	0.40
1:QA:1129:C:H4'	1:QA:1130:A:H5'	2.03	0.40
1:QA:1336:C:H1'	1:QA:1337:G:N1	2.36	0.40
1:QA:1343:G:C6	1:QA:1344:C:N4	2.90	0.40
3:QC:7:PRO:O	3:QC:11:ARG:NH1	2.54	0.40
3:QC:43:LEU:HD22	3:QC:47:LEU:HD22	2.02	0.40
4:QD:135:LEU:HD13	4:QD:135:LEU:HA	1.91	0.40
6:QF:30:LEU:HD23	6:QF:75:LEU:HD11	2.02	0.40
12:QL:22:SER:C	12:QL:24:VAL:N	2.79	0.40
12:QL:53:ARG:HH12	12:QL:92:ASP:HB3	1.86	0.40
19:QS:42:PRO:HG2	50:R4:63:TYR:HE2	1.86	0.40
25:RA:492:A:O5'	25:RA:492:A:H8	2.04	0.40
25:RA:552:G:H2'	25:RA:553:U:O4'	2.22	0.40
25:RA:827:U:H1'	25:RA:2246:G:HO2'	1.86	0.40
25:RA:1434:A:H61	25:RA:1558:A:N6	2.18	0.40
25:RA:2296:U:OP2	38:RS:6:ALA:HB2	2.21	0.40
26:RB:86:G:H2'	26:RB:87:G:H8	1.86	0.40
27:RD:35:LYS:HB3	27:RD:63:ARG:HA	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:RE:92:THR:HB	28:RE:93:VAL:H	1.57	0.40
28:RE:119:ARG:HG2	28:RE:160:TYR:HB2	2.04	0.40
31:RH:65:HIS:O	31:RH:66:GLY:C	2.63	0.40
34:RO:66:LYS:HA	34:RO:79:PHE:O	2.21	0.40
36:RQ:66:ILE:O	36:RQ:67:ARG:HB2	2.22	0.40
44:RY:11:ASP:O	44:RY:26:LYS:HG3	2.21	0.40
50:R4:49:PHE:O	50:R4:50:VAL:CG2	2.69	0.40
51:R5:6:VAL:HA	51:R5:7:PRO:HD3	1.81	0.40
1:XA:675:A:N6	1:XA:715:A:H61	2.19	0.40
1:XA:720:C:H5"	18:XR:52:PRO:HA	2.02	0.40
1:XA:950:U:H2'	1:XA:951:G:C8	2.56	0.40
1:XA:958:A:N6	1:XA:959:A:N1	2.70	0.40
1:XA:1072:G:C5	1:XA:1073:U:C4	3.09	0.40
1:XA:1191:A:H5"	3:XC:4:LYS:NZ	2.34	0.40
2:XB:15:VAL:C	2:XB:16:HIS:CG	2.99	0.40
2:XB:19:HIS:NE2	2:XB:206:ASP:HB2	2.36	0.40
2:XB:74:LYS:HE2	2:XB:74:LYS:HB3	1.93	0.40
5:XE:34:VAL:HG11	5:XE:63:ARG:HG2	2.03	0.40
11:XK:18:ARG:HB3	11:XK:33:THR:OG1	2.21	0.40
11:XK:38:ASN:HA	11:XK:39:PRO:HD3	1.75	0.40
19:XS:15:LEU:HA	19:XS:18:LYS:HB3	2.04	0.40
25:YA:988:A:O5'	25:YA:988:A:H8	2.04	0.40
25:YA:1054:A:H5'	25:YA:1055:G:OP2	2.21	0.40
25:YA:1667:G:O2'	25:YA:1991:U:O4	2.24	0.40
25:YA:1790:C:H2'	25:YA:1791:A:C5	2.56	0.40
25:YA:2626:C:H2'	25:YA:2627:G:H8	1.86	0.40
25:YA:2649:U:H2'	25:YA:2650:U:C6	2.56	0.40
25:YA:2776:A:OP1	25:YA:2776:A:H3'	2.22	0.40
27:YD:72:LYS:HG2	27:YD:103:ARG:HH22	1.85	0.40
27:YD:177:LEU:C	27:YD:179:SER:N	2.80	0.40
28:YE:93:VAL:H	28:YE:95:ILE:CD1	2.22	0.40
28:YE:119:ARG:HG2	28:YE:160:TYR:HB2	2.04	0.40
29:YF:59:TYR:HB3	29:YF:60:SER:H	1.70	0.40
29:YF:118:ALA:C	29:YF:120:GLU:N	2.78	0.40
30:YG:61:ALA:HA	30:YG:64:THR:HG22	2.02	0.40
36:YQ:20:ALA:HB1	36:YQ:99:PRO:CG	2.51	0.40
38:YS:99:LYS:HE2	38:YS:103:GLU:OE2	2.20	0.40
40:YU:17:ILE:HD13	40:YU:17:ILE:HA	1.92	0.40
40:YU:19:LYS:O	40:YU:22:LYS:HB2	2.22	0.40
49:Y3:52:HIS:CD2	49:Y3:53:LEU:HG	2.57	0.40
1:QA:222:U:H2'	1:QA:223:U:H6	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:QA:358:U:H2'	1:QA:359:U:C6	2.56	0.40
1:QA:370:C:H2'	1:QA:371:G:C8	2.56	0.40
1:QA:440:A:H8	1:QA:440:A:OP2	2.04	0.40
1:QA:864:A:H5'	5:QE:86:ALA:HB2	2.03	0.40
1:QA:975:A:H8	1:QA:975:A:H5'	1.85	0.40
1:QA:1235:U:O2'	1:QA:1305:G:O5'	2.39	0.40
1:QA:1392:G:N2	1:QA:1502:A:C8	2.89	0.40
4:QD:63:LYS:HB2	4:QD:63:LYS:HE3	1.78	0.40
9:QI:40:LEU:C	9:QI:42:ARG:H	2.30	0.40
11:QK:74:ALA:C	11:QK:76:GLY:H	2.30	0.40
13:QM:121:LYS:C	13:QM:122:LYS:HG3	2.47	0.40
25:RA:1453:A:C5	25:RA:2702:U:H6	2.40	0.40
25:RA:1518:C:H2'	25:RA:1519:G:C8	2.57	0.40
25:RA:2410:G:H2'	25:RA:2411:A:C8	2.57	0.40
25:RA:2626:C:H2'	25:RA:2627:G:H8	1.86	0.40
27:RD:209:ALA:O	27:RD:212:SER:HB2	2.22	0.40
28:RE:5:LEU:O	28:RE:28:ALA:HA	2.22	0.40
28:RE:27:LEU:HD21	39:RT:1:MET:CE	2.39	0.40
29:RF:7:TYR:O	29:RF:21:ALA:HA	2.21	0.40
33:RN:104:LYS:C	33:RN:106:MET:H	2.28	0.40
36:RQ:46:GLN:OE1	36:RQ:126:PRO:HG3	2.22	0.40
38:RS:69:VAL:HG13	38:RS:101:LEU:HD22	2.03	0.40
42:RW:38:TYR:CD2	51:R5:30:LEU:HD21	2.57	0.40
45:RZ:103:ARG:HD3	45:RZ:136:PHE:CD2	2.55	0.40
50:R4:21:VAL:O	50:R4:22:ILE:O	2.40	0.40
50:R4:26:SER:O	50:R4:27:THR:O	2.40	0.40
53:R7:31:LEU:HD23	53:R7:31:LEU:HA	1.86	0.40
54:R8:53:PRO:HD2	54:R8:54:GLU:H	1.84	0.40
54:R8:53:PRO:CG	54:R8:54:GLU:N	2.84	0.40
1:XA:564:C:H5"	12:XL:10:LEU:CD1	2.51	0.40
1:XA:942:G:C2	1:XA:1342:C:C2	3.10	0.40
1:XA:1461:G:H2'	1:XA:1462:G:H8	1.87	0.40
10:XJ:55:LYS:HE3	10:XJ:56:HIS:CD2	2.55	0.40
12:XL:43:VAL:HG13	12:XL:55:VAL:HG21	2.03	0.40
25:YA:569:U:C4	25:YA:570:G:C6	3.10	0.40
25:YA:924:C:H2'	25:YA:925:C:C6	2.57	0.40
25:YA:1019:U:O2'	25:YA:1021:A:C2	2.75	0.40
25:YA:1320:C:C5	25:YA:1329:U:H5'	2.56	0.40
25:YA:2061:G:H2'	25:YA:2501:C:O2'	2.21	0.40
25:YA:2468:G:H5"	36:YQ:120:ILE:CD1	2.51	0.40
27:YD:13:ARG:HG2	27:YD:13:ARG:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:YD:230:ASP:OD1	27:YD:230:ASP:N	2.54	0.40
28:YE:93:VAL:HG21	28:YE:180:ASN:HA	2.03	0.40
29:YF:33:LEU:O	29:YF:37:VAL:HG23	2.21	0.40
29:YF:36:VAL:HG11	29:YF:183:VAL:HG11	2.04	0.40
29:YF:127:GLU:C	29:YF:129:PHE:H	2.26	0.40
31:YH:128:PRO:CG	31:YH:129:THR:H	2.33	0.40
32:YI:75:LEU:HD23	32:YI:105:HIS:HD2	1.85	0.40
32:YI:144:VAL:O	32:YI:145:VAL:HG22	2.20	0.40
36:YQ:39:PRO:HA	36:YQ:97:VAL:O	2.21	0.40
36:YQ:139:GLU:HG2	36:YQ:140:ALA:N	2.36	0.40
38:YS:24:LEU:HD22	38:YS:24:LEU:N	2.37	0.40
40:YU:61:TRP:O	40:YU:65:ILE:HG13	2.22	0.40
45:YZ:45:ASP:O	45:YZ:49:ARG:HG2	2.22	0.40
50:Y4:14:ILE:O	50:Y4:14:ILE:HG23	2.21	0.40
54:Y8:39:LYS:HD2	54:Y8:39:LYS:O	2.22	0.40
1:QA:224:C:H2'	1:QA:225:C:H6	1.86	0.40
1:QA:1176:A:H2'	1:QA:1177:G:H5'	2.03	0.40
1:QA:1417:G:C6	1:QA:1482:G:C6	3.10	0.40
5:QE:50:GLU:HG3	5:QE:52:PRO:HD2	2.04	0.40
6:QF:21:LEU:O	6:QF:25:ILE:HG12	2.21	0.40
12:QL:11:VAL:HG11	17:QQ:36:ILE:HG21	2.03	0.40
13:QM:82:MET:HE3	13:QM:93:ARG:HG2	2.04	0.40
15:QO:32:LEU:HD23	15:QO:32:LEU:HA	1.76	0.40
25:RA:443:A:H1'	25:RA:1201:C:O4'	2.21	0.40
25:RA:483:A:H1'	44:RY:59:GLY:O	2.21	0.40
25:RA:581:C:OP1	40:RU:31:SER:OG	2.31	0.40
25:RA:729:G:C8	27:RD:208:LYS:HD2	2.56	0.40
25:RA:1341:U:OP2	25:RA:1394:U:O2'	2.25	0.40
25:RA:1359:A:C6	25:RA:1373:A:C5	3.09	0.40
25:RA:2038:G:H2'	25:RA:2039:C:O4'	2.21	0.40
25:RA:2413:G:H21	35:RP:70:GLN:NE2	2.19	0.40
25:RA:2626:C:H2'	25:RA:2627:G:C8	2.56	0.40
26:RB:61:G:C6	26:RB:62:C:C4	3.10	0.40
30:RG:117:PHE:HE1	30:RG:120:LEU:HD23	1.87	0.40
32:RI:88:ILE:HG12	32:RI:122:GLU:N	2.36	0.40
33:RN:17:ASP:O	33:RN:19:GLU:N	2.54	0.40
34:RO:68:GLU:H	34:RO:68:GLU:CD	2.29	0.40
35:RP:18:ARG:HH11	35:RP:27:HIS:CD2	2.40	0.40
36:RQ:51:ARG:O	36:RQ:52:VAL:C	2.62	0.40
36:RQ:119:ARG:NH1	36:RQ:119:ARG:CG	2.83	0.40
44:RY:67:LEU:HD12	44:RY:67:LEU:HA	1.93	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:R0:68:GLU:HG2	46:R0:80:HIS:HB2	2.04	0.40
47:R1:87:PRO:O	47:R1:91:LYS:HB2	2.21	0.40
50:R4:7:PRO:O	50:R4:8:LYS:C	2.64	0.40
54:R8:40:GLU:O	54:R8:42:ARG:N	2.54	0.40
1:XA:129(A):G:N2	1:XA:191(A):G:C5	2.89	0.40
1:XA:384:G:H2'	1:XA:385:C:C6	2.56	0.40
1:XA:1051:C:H2'	1:XA:1052:U:H6	1.87	0.40
1:XA:1082:G:H2'	1:XA:1083:U:O4'	2.22	0.40
1:XA:1292:U:OP2	7:XG:41:ARG:NH2	2.55	0.40
2:XB:166:ASP:HB3	2:XB:169:LYS:HB2	2.03	0.40
5:XE:89:ILE:HG12	5:XE:91:LEU:CD1	2.52	0.40
9:XI:5:TYR:HA	9:XI:17:VAL:O	2.21	0.40
10:XJ:77:PRO:O	10:XJ:79:ARG:NH1	2.54	0.40
17:XQ:58:GLU:O	17:XQ:74:LEU:N	2.40	0.40
25:YA:896:A:C5	45:YZ:146:ILE:HD12	2.57	0.40
25:YA:1050:A:C8	25:YA:2751:G:C4	3.09	0.40
25:YA:1473:G:H1	25:YA:1520:U:H3	1.69	0.40
25:YA:2079:U:H2'	25:YA:2080:G:O4'	2.21	0.40
25:YA:2344:U:O2'	52:Y6:37:ARG:HG2	2.21	0.40
25:YA:2461:C:H2'	25:YA:2462:U:C6	2.57	0.40
25:YA:2483:C:O2'	36:YQ:49:ALA:CB	2.70	0.40
25:YA:2745:C:H1'	31:YH:143:GLN:HG2	2.03	0.40
25:YA:2811:G:OP1	28:YE:61:ARG:HG2	2.20	0.40
27:YD:142:VAL:HA	27:YD:194:GLY:H	1.86	0.40
28:YE:94:GLU:C	28:YE:96:PHE:H	2.29	0.40
29:YF:124:LEU:HD12	29:YF:125:LEU:O	2.22	0.40
29:YF:160:ASN:O	29:YF:161:GLU:C	2.64	0.40
41:YV:3:ALA:HA	41:YV:40:LEU:O	2.21	0.40
41:YV:21:ARG:HD2	41:YV:91:TYR:CE1	2.56	0.40
52:Y6:28:ARG:HA	52:Y6:28:ARG:HD2	1.86	0.40
54:Y8:53:PRO:O	54:Y8:54:GLU:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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	QB	235/256 (92%)	173 (74%)	46 (20%)	16 (7%)	1 11
2	XB	235/256 (92%)	178 (76%)	42 (18%)	15 (6%)	1 12
3	QC	203/239 (85%)	163 (80%)	34 (17%)	6 (3%)	3 25
3	XC	203/239 (85%)	172 (85%)	28 (14%)	3 (2%)	8 37
4	QD	206/209 (99%)	186 (90%)	17 (8%)	3 (2%)	8 37
4	XD	206/209 (99%)	176 (85%)	25 (12%)	5 (2%)	4 29
5	QE	149/162 (92%)	137 (92%)	8 (5%)	4 (3%)	4 27
5	XE	149/162 (92%)	134 (90%)	12 (8%)	3 (2%)	6 32
6	QF	99/101 (98%)	95 (96%)	4 (4%)	0	100 100
6	XF	99/101 (98%)	94 (95%)	5 (5%)	0	100 100
7	QG	153/156 (98%)	135 (88%)	16 (10%)	2 (1%)	9 39
7	XG	153/156 (98%)	138 (90%)	13 (8%)	2 (1%)	9 39
8	QH	136/138 (99%)	121 (89%)	14 (10%)	1 (1%)	18 51
8	XH	136/138 (99%)	120 (88%)	12 (9%)	4 (3%)	3 26
9	QI	125/128 (98%)	103 (82%)	17 (14%)	5 (4%)	2 20
9	XI	125/128 (98%)	97 (78%)	24 (19%)	4 (3%)	3 24
10	QJ	97/105 (92%)	77 (79%)	16 (16%)	4 (4%)	2 19
10	XJ	97/105 (92%)	79 (81%)	13 (13%)	5 (5%)	1 15
11	QK	117/129 (91%)	101 (86%)	14 (12%)	2 (2%)	7 35
11	XK	117/129 (91%)	101 (86%)	14 (12%)	2 (2%)	7 35
12	QL	123/132 (93%)	85 (69%)	24 (20%)	14 (11%)	0 4
12	XL	123/132 (93%)	85 (69%)	24 (20%)	14 (11%)	0 4
13	QM	119/126 (94%)	95 (80%)	15 (13%)	9 (8%)	1 9
13	XM	119/126 (94%)	94 (79%)	16 (13%)	9 (8%)	1 9
14	QN	58/61 (95%)	50 (86%)	4 (7%)	4 (7%)	1 11
14	XN	58/61 (95%)	46 (79%)	6 (10%)	6 (10%)	0 5
15	QO	86/89 (97%)	80 (93%)	5 (6%)	1 (1%)	10 41
15	XO	86/89 (97%)	80 (93%)	4 (5%)	2 (2%)	5 30
16	QP	82/88 (93%)	73 (89%)	8 (10%)	1 (1%)	10 41
16	XP	82/88 (93%)	72 (88%)	9 (11%)	1 (1%)	10 41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
17	QQ	98/105 (93%)	91 (93%)	5 (5%)	2 (2%)	6 32
17	XQ	98/105 (93%)	88 (90%)	10 (10%)	0	100 100
18	QR	68/88 (77%)	56 (82%)	9 (13%)	3 (4%)	2 18
18	XR	68/88 (77%)	61 (90%)	6 (9%)	1 (2%)	8 37
19	QS	82/93 (88%)	55 (67%)	16 (20%)	11 (13%)	0 3
19	XS	82/93 (88%)	54 (66%)	18 (22%)	10 (12%)	0 4
20	QT	97/106 (92%)	76 (78%)	15 (16%)	6 (6%)	1 12
20	XT	97/106 (92%)	75 (77%)	16 (16%)	6 (6%)	1 12
21	QU	23/27 (85%)	19 (83%)	3 (13%)	1 (4%)	2 18
21	XU	23/27 (85%)	18 (78%)	4 (17%)	1 (4%)	2 18
27	RD	270/276 (98%)	226 (84%)	32 (12%)	12 (4%)	2 18
27	YD	270/276 (98%)	204 (76%)	47 (17%)	19 (7%)	1 10
28	RE	203/206 (98%)	120 (59%)	41 (20%)	42 (21%)	0 1
28	YE	203/206 (98%)	120 (59%)	41 (20%)	42 (21%)	0 1
29	RF	200/210 (95%)	181 (90%)	13 (6%)	6 (3%)	3 25
29	YF	200/210 (95%)	143 (72%)	37 (18%)	20 (10%)	0 5
30	RG	179/182 (98%)	139 (78%)	25 (14%)	15 (8%)	0 7
30	YG	179/182 (98%)	142 (79%)	25 (14%)	12 (7%)	1 11
31	RH	168/180 (93%)	94 (56%)	36 (21%)	38 (23%)	0 0
31	YH	168/180 (93%)	94 (56%)	36 (21%)	38 (23%)	0 0
32	RI	144/148 (97%)	109 (76%)	24 (17%)	11 (8%)	1 9
32	YI	144/148 (97%)	108 (75%)	23 (16%)	13 (9%)	0 6
33	RN	136/140 (97%)	104 (76%)	20 (15%)	12 (9%)	0 7
33	YN	136/140 (97%)	107 (79%)	16 (12%)	13 (10%)	0 6
34	RO	120/122 (98%)	109 (91%)	9 (8%)	2 (2%)	7 35
34	YO	120/122 (98%)	108 (90%)	10 (8%)	2 (2%)	7 35
35	RP	148/150 (99%)	109 (74%)	26 (18%)	13 (9%)	0 7
35	YP	148/150 (99%)	101 (68%)	35 (24%)	12 (8%)	1 8
36	RQ	139/141 (99%)	95 (68%)	30 (22%)	14 (10%)	0 5
36	YQ	139/141 (99%)	97 (70%)	28 (20%)	14 (10%)	0 5
37	RR	116/118 (98%)	106 (91%)	5 (4%)	5 (4%)	2 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
37	YR	116/118 (98%)	99 (85%)	11 (10%)	6 (5%)	1 15
38	RS	109/112 (97%)	76 (70%)	22 (20%)	11 (10%)	0 5
38	YS	109/112 (97%)	62 (57%)	29 (27%)	18 (16%)	0 2
39	RT	135/146 (92%)	107 (79%)	16 (12%)	12 (9%)	0 7
39	YT	135/146 (92%)	108 (80%)	17 (13%)	10 (7%)	1 9
40	RU	115/118 (98%)	102 (89%)	9 (8%)	4 (4%)	3 23
40	YU	115/118 (98%)	102 (89%)	11 (10%)	2 (2%)	7 35
41	RV	99/101 (98%)	82 (83%)	11 (11%)	6 (6%)	1 13
41	YV	99/101 (98%)	79 (80%)	12 (12%)	8 (8%)	1 8
42	RW	111/113 (98%)	99 (89%)	8 (7%)	4 (4%)	2 22
42	YW	111/113 (98%)	100 (90%)	9 (8%)	2 (2%)	6 34
43	RX	90/96 (94%)	77 (86%)	11 (12%)	2 (2%)	5 30
43	YX	90/96 (94%)	82 (91%)	6 (7%)	2 (2%)	5 30
44	RY	100/110 (91%)	71 (71%)	13 (13%)	16 (16%)	0 2
44	YY	100/110 (91%)	70 (70%)	18 (18%)	12 (12%)	0 4
45	RZ	181/206 (88%)	127 (70%)	34 (19%)	20 (11%)	0 5
45	YZ	181/206 (88%)	135 (75%)	28 (16%)	18 (10%)	0 6
46	R0	80/85 (94%)	66 (82%)	13 (16%)	1 (1%)	9 39
46	Y0	80/85 (94%)	73 (91%)	7 (9%)	0	100 100
47	R1	95/98 (97%)	75 (79%)	11 (12%)	9 (10%)	0 6
47	Y1	95/98 (97%)	72 (76%)	17 (18%)	6 (6%)	1 12
48	R2	67/72 (93%)	54 (81%)	9 (13%)	4 (6%)	1 13
48	Y2	67/72 (93%)	46 (69%)	13 (19%)	8 (12%)	0 4
49	R3	57/60 (95%)	52 (91%)	3 (5%)	2 (4%)	3 23
49	Y3	57/60 (95%)	52 (91%)	4 (7%)	1 (2%)	6 34
50	R4	69/71 (97%)	22 (32%)	21 (30%)	26 (38%)	0 0
50	Y4	69/71 (97%)	35 (51%)	15 (22%)	19 (28%)	0 0
51	R5	57/60 (95%)	33 (58%)	9 (16%)	15 (26%)	0 0
51	Y5	56/60 (93%)	46 (82%)	8 (14%)	2 (4%)	2 22
52	R6	47/54 (87%)	23 (49%)	13 (28%)	11 (23%)	0 0
52	Y6	47/54 (87%)	22 (47%)	17 (36%)	8 (17%)	0 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
53	R7	47/49 (96%)	45 (96%)	1 (2%)	1 (2%)	5 31
53	Y7	47/49 (96%)	43 (92%)	3 (6%)	1 (2%)	5 31
54	R8	62/65 (95%)	36 (58%)	14 (23%)	12 (19%)	0 1
54	Y8	62/65 (95%)	36 (58%)	15 (24%)	11 (18%)	0 1
55	R9	35/37 (95%)	35 (100%)	0	0	100 100
55	Y9	35/37 (95%)	31 (89%)	4 (11%)	0	100 100
All	All	11469/12128 (95%)	9024 (79%)	1612 (14%)	833 (7%)	1 10

All (833) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	QB	236	TYR
3	QC	12	LEU
3	QC	190	ARG
4	QD	33	MET
12	QL	18	VAL
12	QL	27	LEU
12	QL	62	SER
12	QL	121	GLY
13	QM	67	GLU
13	QM	106	ASN
13	QM	118	ALA
14	QN	16	PHE
19	QS	12	ASP
19	QS	45	VAL
20	QT	49	ALA
27	RD	26	LYS
27	RD	122	ASP
27	RD	242	ARG
28	RE	4	ILE
28	RE	7	VAL
28	RE	9	VAL
28	RE	22	PRO
28	RE	54	GLN
28	RE	57	LYS
28	RE	60	ASN
28	RE	63	LEU
28	RE	64	LYS
28	RE	68	ALA
28	RE	70	ALA

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Mol	Chain	Res	Type
28	RE	73	GLU
28	RE	90	THR
28	RE	92	THR
28	RE	93	VAL
28	RE	169	ASN
28	RE	187	ALA
28	RE	189	PRO
29	RF	89	VAL
29	RF	134	GLY
31	RH	10	PRO
31	RH	12	PRO
31	RH	83	TYR
31	RH	85	LYS
31	RH	86	GLU
31	RH	87	LEU
31	RH	90	LYS
31	RH	92	ILE
31	RH	126	PRO
31	RH	127	GLU
31	RH	128	PRO
31	RH	137	ASP
31	RH	138	LYS
31	RH	153	LYS
31	RH	154	PRO
31	RH	155	SER
31	RH	169	VAL
32	RI	115	ALA
32	RI	145	VAL
33	RN	9	VAL
33	RN	22	THR
33	RN	96	GLU
33	RN	131	GLN
34	RO	5	GLN
35	RP	15	ARG
35	RP	61	ARG
35	RP	148	LEU
36	RQ	6	ARG
36	RQ	18	LYS
36	RQ	22	LYS
36	RQ	27	VAL
36	RQ	81	VAL
36	RQ	90	VAL

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Mol	Chain	Res	Type
36	RQ	134	ARG
37	RR	3	HIS
37	RR	4	LEU
38	RS	57	LYS
38	RS	88	ASP
38	RS	89	ARG
39	RT	2	ASN
39	RT	3	ARG
39	RT	106	SER
39	RT	112	ARG
39	RT	124	ASP
40	RU	91	ASP
41	RV	48	GLY
41	RV	50	PRO
41	RV	100	ARG
42	RW	111	HIS
44	RY	3	VAL
44	RY	50	ARG
44	RY	57	GLN
44	RY	77	PRO
44	RY	78	ALA
45	RZ	6	LYS
45	RZ	111	VAL
48	R2	70	GLN
48	R2	71	ASN
50	R4	5	ILE
50	R4	14	ILE
50	R4	16	CYS
50	R4	22	ILE
50	R4	23	GLU
50	R4	36	CYS
50	R4	37	SER
50	R4	40	HIS
50	R4	42	PHE
50	R4	43	TYR
50	R4	49	PHE
50	R4	50	VAL
50	R4	51	ASP
50	R4	53	GLU
50	R4	62	ARG
50	R4	66	SER
50	R4	68	ARG

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Mol	Chain	Res	Type
51	R5	3	LYS
51	R5	4	HIS
51	R5	35	GLU
51	R5	51	TYR
51	R5	53	ALA
52	R6	15	GLU
54	R8	29	LYS
54	R8	31	HIS
54	R8	34	TRP
54	R8	52	LYS
54	R8	62	LEU
2	XB	230	VAL
2	XB	236	TYR
3	XC	12	LEU
3	XC	79	ARG
4	XD	154	ASN
11	XK	91	ARG
12	XL	18	VAL
12	XL	27	LEU
12	XL	62	SER
12	XL	121	GLY
13	XM	67	GLU
13	XM	106	ASN
13	XM	118	ALA
14	XN	14	PRO
14	XN	16	PHE
14	XN	52	GLN
19	XS	3	ARG
19	XS	12	ASP
20	XT	96	GLY
27	YD	26	LYS
27	YD	28	GLU
27	YD	123	ALA
27	YD	231	HIS
28	YE	4	ILE
28	YE	7	VAL
28	YE	9	VAL
28	YE	22	PRO
28	YE	54	GLN
28	YE	57	LYS
28	YE	60	ASN
28	YE	63	LEU

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Mol	Chain	Res	Type
28	YE	64	LYS
28	YE	68	ALA
28	YE	70	ALA
28	YE	73	GLU
28	YE	90	THR
28	YE	92	THR
28	YE	93	VAL
28	YE	169	ASN
28	YE	187	ALA
28	YE	189	PRO
29	YF	25	PRO
29	YF	66	PRO
29	YF	68	LYS
29	YF	73	ALA
29	YF	89	VAL
29	YF	128	ALA
29	YF	176	LEU
30	YG	96	ARG
31	YH	10	PRO
31	YH	12	PRO
31	YH	83	TYR
31	YH	85	LYS
31	YH	86	GLU
31	YH	87	LEU
31	YH	90	LYS
31	YH	92	ILE
31	YH	126	PRO
31	YH	127	GLU
31	YH	128	PRO
31	YH	137	ASP
31	YH	138	LYS
31	YH	153	LYS
31	YH	154	PRO
31	YH	155	SER
31	YH	169	VAL
32	YI	133	HIS
32	YI	145	VAL
33	YN	9	VAL
33	YN	22	THR
33	YN	36	GLY
35	YP	6	LEU
35	YP	15	ARG

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Mol	Chain	Res	Type
35	YP	95	VAL
35	YP	148	LEU
36	YQ	6	ARG
36	YQ	18	LYS
36	YQ	22	LYS
36	YQ	27	VAL
36	YQ	81	VAL
36	YQ	90	VAL
36	YQ	134	ARG
37	YR	3	HIS
38	YS	4	LEU
38	YS	12	PHE
38	YS	14	VAL
38	YS	23	ARG
38	YS	56	LEU
38	YS	57	LYS
38	YS	88	ASP
38	YS	89	ARG
38	YS	90	GLY
38	YS	107	GLU
39	YT	2	ASN
39	YT	123	GLN
39	YT	124	ASP
40	YU	93	LYS
41	YV	45	THR
43	YX	68	ARG
44	YY	50	ARG
44	YY	57	GLN
44	YY	77	PRO
44	YY	78	ALA
45	YZ	6	LYS
45	YZ	146	ILE
45	YZ	152	ALA
45	YZ	159	PRO
45	YZ	166	SER
47	Y1	30	VAL
47	Y1	84	GLY
47	Y1	91	LYS
47	Y1	95	LEU
48	Y2	16	LEU
48	Y2	43	GLN
48	Y2	47	ASN

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Mol	Chain	Res	Type
48	Y2	48	HIS
48	Y2	71	ASN
50	Y4	24	THR
50	Y4	40	HIS
50	Y4	49	PHE
52	Y6	15	GLU
53	Y7	48	LYS
54	Y8	29	LYS
54	Y8	31	HIS
54	Y8	34	TRP
54	Y8	52	LYS
54	Y8	62	LEU
2	QB	15	VAL
2	QB	96	ARG
2	QB	229	VAL
2	QB	230	VAL
2	QB	237	ALA
3	QC	79	ARG
4	QD	166	LYS
5	QE	115	VAL
8	QH	129	VAL
9	QI	41	VAL
9	QI	117	HIS
10	QJ	59	SER
11	QK	103	LEU
12	QL	49	ASN
12	QL	65	GLU
12	QL	110	VAL
12	QL	115	LYS
12	QL	116	SER
12	QL	128	ALA
13	QM	12	ASN
13	QM	120	LYS
14	QN	12	ARG
17	QQ	81	ARG
19	QS	3	ARG
19	QS	11	VAL
19	QS	26	GLY
19	QS	31	ILE
19	QS	41	VAL
27	RD	32	SER
28	RE	8	LYS

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Mol	Chain	Res	Type
28	RE	37	ARG
28	RE	53	PRO
28	RE	61	ARG
28	RE	78	LEU
28	RE	88	GLY
28	RE	186	GLY
28	RE	190	GLY
28	RE	204	ALA
29	RF	67	GLN
29	RF	73	ALA
30	RG	4	ASP
30	RG	5	VAL
30	RG	14	GLU
30	RG	96	ARG
30	RG	137	GLU
30	RG	146	TYR
31	RH	3	ARG
31	RH	8	PRO
31	RH	55	PRO
31	RH	59	ARG
31	RH	84	SER
31	RH	151	ILE
31	RH	156	ALA
31	RH	168	PRO
32	RI	10	GLU
32	RI	11	ASN
32	RI	15	VAL
32	RI	133	HIS
35	RP	6	LEU
35	RP	65	ARG
35	RP	67	MET
35	RP	103	ALA
35	RP	106	LEU
35	RP	141	ALA
36	RQ	13	GLN
36	RQ	24	GLY
36	RQ	28	ALA
36	RQ	57	HIS
37	RR	107	ASP
38	RS	4	LEU
38	RS	107	GLU
39	RT	37	GLY

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Mol	Chain	Res	Type
40	RU	90	VAL
41	RV	49	THR
41	RV	79	VAL
43	RX	41	ASN
44	RY	45	VAL
44	RY	48	ALA
44	RY	63	LYS
45	RZ	51	ALA
45	RZ	108	PRO
45	RZ	116	VAL
45	RZ	153	SER
45	RZ	177	PRO
47	R1	30	VAL
47	R1	80	LEU
47	R1	84	GLY
47	R1	91	LYS
47	R1	95	LEU
48	R2	43	GLN
49	R3	26	LEU
50	R4	9	LEU
50	R4	24	THR
51	R5	43	HIS
51	R5	55	ARG
52	R6	7	ILE
52	R6	45	LYS
2	XB	15	VAL
4	XD	30	LYS
4	XD	166	LYS
5	XE	115	VAL
7	XG	55	GLY
8	XH	50	ARG
9	XI	41	VAL
9	XI	127	LYS
10	XJ	30	SER
10	XJ	86	MET
12	XL	65	GLU
12	XL	110	VAL
12	XL	115	LYS
12	XL	116	SER
12	XL	128	ALA
13	XM	6	GLY
13	XM	21	TYR

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Mol	Chain	Res	Type
19	XS	41	VAL
19	XS	45	VAL
20	XT	99	LEU
27	YD	3	VAL
27	YD	32	SER
27	YD	58	HIS
27	YD	122	ASP
27	YD	169	GLU
28	YE	8	LYS
28	YE	20	ALA
28	YE	53	PRO
28	YE	61	ARG
28	YE	78	LEU
28	YE	88	GLY
28	YE	186	GLY
28	YE	190	GLY
28	YE	204	ALA
29	YF	18	ARG
29	YF	107	LYS
29	YF	108	LYS
29	YF	111	ALA
29	YF	132	VAL
29	YF	134	GLY
29	YF	168	ARG
30	YG	4	ASP
30	YG	36	LYS
31	YH	3	ARG
31	YH	8	PRO
31	YH	55	PRO
31	YH	59	ARG
31	YH	84	SER
31	YH	151	ILE
31	YH	156	ALA
31	YH	168	PRO
32	YI	12	LEU
33	YN	23	LEU
33	YN	96	GLU
34	YO	5	GLN
35	YP	106	LEU
36	YQ	13	GLN
36	YQ	24	GLY
36	YQ	28	ALA

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Mol	Chain	Res	Type
37	YR	4	LEU
37	YR	45	ARG
37	YR	107	ASP
38	YS	61	ASN
38	YS	87	PHE
38	YS	96	GLY
38	YS	100	ALA
38	YS	109	GLY
39	YT	13	ARG
39	YT	39	ARG
39	YT	106	SER
41	YV	31	ALA
41	YV	48	GLY
41	YV	79	VAL
42	YW	111	HIS
44	YY	58	GLY
44	YY	102	CYS
45	YZ	53	ILE
45	YZ	59	LEU
45	YZ	81	ARG
48	Y2	24	LEU
48	Y2	44	LEU
48	Y2	70	GLN
50	Y4	5	ILE
50	Y4	18	CYS
50	Y4	22	ILE
50	Y4	37	SER
50	Y4	43	TYR
50	Y4	50	VAL
52	Y6	7	ILE
52	Y6	16	CYS
52	Y6	33	LYS
2	QB	26	PRO
2	QB	87	ARG
2	QB	207	ALA
3	QC	4	LYS
3	QC	51	GLY
4	QD	155	LEU
5	QE	77	PRO
10	QJ	30	SER
11	QK	125	PHE
12	QL	51	ALA

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Mol	Chain	Res	Type
12	QL	123	LYS
14	QN	14	PRO
15	QO	23	GLY
17	QQ	74	LEU
19	QS	9	VAL
19	QS	14	HIS
19	QS	28	LYS
20	QT	96	GLY
27	RD	46	GLN
27	RD	239	ARG
28	RE	20	ALA
28	RE	62	PRO
28	RE	69	LYS
28	RE	71	GLY
28	RE	82	ARG
28	RE	117	MET
28	RE	130	GLY
28	RE	132	HIS
29	RF	66	PRO
29	RF	133	ASN
30	RG	32	PRO
30	RG	116	ASP
31	RH	50	VAL
31	RH	81	GLU
31	RH	152	ARG
32	RI	102	SER
32	RI	118	LYS
32	RI	122	GLU
33	RN	8	GLN
33	RN	23	LEU
33	RN	95	PRO
33	RN	130	HIS
34	RO	97	ARG
35	RP	90	ARG
36	RQ	88	GLY
36	RQ	91	GLU
37	RR	74	LYS
38	RS	12	PHE
38	RS	61	ASN
39	RT	12	SER
39	RT	97	ALA
40	RU	117	GLN

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Mol	Chain	Res	Type
43	RX	67	GLY
44	RY	58	GLY
45	RZ	166	SER
47	R1	76	ARG
48	R2	47	ASN
49	R3	27	GLY
50	R4	27	THR
50	R4	46	GLN
52	R6	33	LYS
52	R6	35	GLU
52	R6	49	HIS
54	R8	46	ARG
54	R8	47	LYS
2	XB	13	ALA
2	XB	22	LYS
2	XB	24	TRP
2	XB	135	GLN
2	XB	207	ALA
4	XD	73	ARG
4	XD	155	LEU
7	XG	7	ALA
8	XH	2	LEU
9	XI	56	LEU
9	XI	95	LYS
10	XJ	59	SER
11	XK	103	LEU
12	XL	51	ALA
12	XL	123	LYS
13	XM	4	ILE
13	XM	12	ASN
13	XM	42	ALA
19	XS	27	GLU
19	XS	28	LYS
27	YD	111	LEU
27	YD	239	ARG
27	YD	242	ARG
27	YD	262	ARG
28	YE	37	ARG
28	YE	62	PRO
28	YE	69	LYS
28	YE	71	GLY
28	YE	82	ARG

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Mol	Chain	Res	Type
28	YE	117	MET
28	YE	130	GLY
28	YE	132	HIS
31	YH	50	VAL
31	YH	81	GLU
31	YH	152	ARG
32	YI	11	ASN
32	YI	113	ARG
32	YI	117	GLU
32	YI	122	GLU
33	YN	131	GLN
35	YP	29	LYS
36	YQ	57	HIS
36	YQ	88	GLY
36	YQ	91	GLU
37	YR	86	ARG
38	YS	19	LYS
38	YS	74	ALA
38	YS	75	GLU
39	YT	97	ALA
41	YV	49	THR
41	YV	53	GLU
41	YV	100	ARG
44	YY	42	VAL
44	YY	63	LYS
45	YZ	13	GLU
45	YZ	92	SER
45	YZ	113	ALA
45	YZ	181	GLU
49	Y3	3	ARG
50	Y4	23	GLU
50	Y4	30	GLU
50	Y4	34	GLU
50	Y4	66	SER
51	Y5	47	PRO
52	Y6	19	ARG
52	Y6	49	HIS
54	Y8	46	ARG
54	Y8	47	LYS
2	QB	22	LYS
2	QB	126	GLU
2	QB	209	ARG

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Mol	Chain	Res	Type
9	QI	56	LEU
12	QL	64	TYR
13	QM	6	GLY
13	QM	13	LYS
18	QR	20	ALA
18	QR	54	ARG
20	QT	71	THR
21	QU	9	ARG
27	RD	3	VAL
27	RD	123	ALA
27	RD	237	GLU
28	RE	66	HIS
28	RE	126	PRO
30	RG	36	LYS
30	RG	86	MET
31	RH	13	LYS
31	RH	109	PHE
31	RH	159	GLU
32	RI	117	GLU
35	RP	21	ARG
35	RP	95	VAL
37	RR	71	GLN
38	RS	109	GLY
39	RT	38	ASN
39	RT	39	ARG
40	RU	98	LEU
42	RW	18	ARG
42	RW	63	ASP
42	RW	68	ARG
44	RY	4	LYS
44	RY	53	PRO
44	RY	99	CYS
45	RZ	7	ALA
45	RZ	13	GLU
45	RZ	92	SER
45	RZ	112	ARG
50	R4	8	LYS
51	R5	14	ALA
51	R5	37	LYS
51	R5	42	PRO
51	R5	45	VAL
51	R5	48	GLU

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Mol	Chain	Res	Type
52	R6	16	CYS
53	R7	48	LYS
54	R8	30	ARG
54	R8	53	PRO
54	R8	57	ARG
2	XB	19	HIS
2	XB	101	MET
2	XB	155	LEU
8	XH	129	VAL
12	XL	64	TYR
14	XN	15	LYS
14	XN	32	SER
15	XO	88	ARG
18	XR	20	ALA
19	XS	9	VAL
20	XT	49	ALA
20	XT	84	LEU
20	XT	98	PRO
21	XU	9	ARG
27	YD	12	SER
27	YD	73	VAL
28	YE	66	HIS
28	YE	126	PRO
29	YF	43	LYS
29	YF	130	ALA
29	YF	145	GLU
30	YG	14	GLU
30	YG	82	LEU
30	YG	86	MET
30	YG	116	ASP
31	YH	13	LYS
31	YH	47	GLU
31	YH	109	PHE
31	YH	159	GLU
32	YI	10	GLU
32	YI	114	LEU
33	YN	11	PRO
33	YN	28	THR
33	YN	47	ALA
35	YP	25	SER
35	YP	61	ARG
35	YP	93	GLY

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Mol	Chain	Res	Type
39	YT	17	THR
43	YX	40	LYS
44	YY	51	VAL
44	YY	53	PRO
50	Y4	9	LEU
50	Y4	16	CYS
50	Y4	25	TYR
50	Y4	54	GLY
50	Y4	60	GLN
52	Y6	35	GLU
54	Y8	25	MET
54	Y8	53	PRO
54	Y8	57	ARG
2	QB	234	PRO
5	QE	96	PRO
7	QG	7	ALA
9	QI	121	ARG
12	QL	63	GLY
13	QM	4	ILE
18	QR	26	LEU
20	QT	73	HIS
20	QT	97	ALA
27	RD	238	GLY
28	RE	79	ARG
30	RG	82	LEU
30	RG	117	PHE
31	RH	11	VAL
31	RH	27	LYS
31	RH	47	GLU
31	RH	77	LYS
31	RH	170	ARG
32	RI	12	LEU
33	RN	18	ALA
33	RN	57	ALA
33	RN	135	PRO
35	RP	57	THR
38	RS	97	ARG
38	RS	110	LEU
39	RT	40	THR
44	RY	39	VAL
44	RY	41	GLY
44	RY	62	GLU

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Mol	Chain	Res	Type
45	RZ	61	LEU
45	RZ	66	SER
45	RZ	81	ARG
45	RZ	130	PRO
46	R0	18	ALA
47	R1	74	VAL
47	R1	82	LEU
50	R4	30	GLU
50	R4	69	LYS
52	R6	9	LEU
52	R6	19	ARG
54	R8	25	MET
54	R8	64	TYR
2	XB	121	LEU
3	XC	181	ASN
5	XE	70	PRO
10	XJ	27	ALA
12	XL	63	GLY
13	XM	101	GLN
15	XO	23	GLY
20	XT	97	ALA
27	YD	33	LEU
28	YE	79	ARG
29	YF	47	GLY
29	YF	118	ALA
29	YF	136	THR
30	YG	5	VAL
30	YG	53	LEU
31	YH	11	VAL
31	YH	27	LYS
31	YH	77	LYS
31	YH	170	ARG
32	YI	18	VAL
32	YI	118	LYS
33	YN	95	PRO
33	YN	127	ASP
33	YN	134	ARG
33	YN	135	PRO
35	YP	7	ARG
35	YP	65	ARG
35	YP	115	LEU
39	YT	86	ILE

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Mol	Chain	Res	Type
40	YU	117	GLN
41	YV	50	PRO
44	YY	39	VAL
45	YZ	7	ALA
45	YZ	61	LEU
45	YZ	143	GLY
45	YZ	153	SER
47	Y1	74	VAL
50	Y4	14	ILE
51	Y5	5	PRO
52	Y6	21	TYR
54	Y8	64	TYR
2	QB	155	LEU
10	QJ	82	ILE
13	QM	10	PRO
14	QN	15	LYS
20	QT	98	PRO
27	RD	125	ILE
30	RG	52	ILE
30	RG	88	ILE
31	RH	7	LEU
31	RH	26	VAL
38	RS	82	ILE
44	RY	5	MET
47	R1	55	GLY
50	R4	33	VAL
50	R4	70	GLY
51	R5	57	VAL
52	R6	21	TYR
52	R6	34	LEU
2	XB	26	PRO
2	XB	126	GLU
2	XB	237	ALA
10	XJ	91	PRO
14	XN	60	SER
27	YD	178	PRO
27	YD	241	PRO
30	YG	117	PHE
31	YH	7	LEU
31	YH	26	VAL
34	YO	97	ARG
44	YY	3	VAL

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Mol	Chain	Res	Type
45	YZ	177	PRO
47	Y1	55	GLY
2	QB	5	ILE
16	QP	46	PRO
28	RE	86	PRO
28	RE	184	VAL
36	RQ	86	GLY
45	RZ	53	ILE
5	XE	74	GLY
28	YE	86	PRO
28	YE	184	VAL
32	YI	15	VAL
36	YQ	86	GLY
3	QC	81	GLY
5	QE	74	GLY
19	QS	46	GLY
27	RD	35	LYS
51	R5	46	CYS
19	XS	26	GLY
19	XS	46	GLY
39	YT	37	GLY
42	YW	14	PRO
45	YZ	160	GLY
7	QG	50	ILE
9	QI	89	ASN
10	QJ	37	PRO
33	RN	134	ARG
39	RT	86	ILE
45	RZ	62	PRO
45	RZ	165	VAL
51	R5	34	PRO
8	XH	51	VAL
16	XP	46	PRO
19	XS	31	ILE
27	YD	34	VAL
28	RE	52	LEU
28	RE	55	ASN
41	RV	54	GLY
28	YE	52	LEU
28	YE	55	ASN
30	YG	52	ILE
30	YG	88	ILE

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Mol	Chain	Res	Type
32	YI	13	GLY
37	YR	117	VAL
2	QB	227	GLY
30	RG	68	PRO
45	RZ	94	GLU
12	XL	48	PRO

5.3.2 Protein sidechains (i)

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	QB	205/220 (93%)	170 (83%)	35 (17%)	2 13
2	XB	205/220 (93%)	180 (88%)	25 (12%)	5 23
3	QC	159/188 (85%)	146 (92%)	13 (8%)	10 36
3	XC	159/188 (85%)	145 (91%)	14 (9%)	9 33
4	QD	180/181 (99%)	161 (89%)	19 (11%)	6 28
4	XD	180/181 (99%)	155 (86%)	25 (14%)	3 19
5	QE	116/123 (94%)	100 (86%)	16 (14%)	3 20
5	XE	116/123 (94%)	100 (86%)	16 (14%)	3 20
6	QF	90/90 (100%)	79 (88%)	11 (12%)	5 23
6	XF	90/90 (100%)	78 (87%)	12 (13%)	4 21
7	QG	126/127 (99%)	115 (91%)	11 (9%)	9 34
7	XG	126/127 (99%)	112 (89%)	14 (11%)	6 27
8	QH	119/119 (100%)	105 (88%)	14 (12%)	5 24
8	XH	119/119 (100%)	106 (89%)	13 (11%)	6 27
9	QI	98/99 (99%)	80 (82%)	18 (18%)	1 10
9	XI	98/99 (99%)	77 (79%)	21 (21%)	1 7
10	QJ	89/92 (97%)	77 (86%)	12 (14%)	4 21
10	XJ	89/92 (97%)	74 (83%)	15 (17%)	2 13
11	QK	90/99 (91%)	79 (88%)	11 (12%)	5 23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
11	XK	90/99 (91%)	82 (91%)	8 (9%)	9 33
12	QL	104/109 (95%)	87 (84%)	17 (16%)	2 15
12	XL	104/109 (95%)	86 (83%)	18 (17%)	2 12
13	QM	97/101 (96%)	73 (75%)	24 (25%)	1 4
13	XM	97/101 (96%)	79 (81%)	18 (19%)	1 10
14	QN	49/50 (98%)	40 (82%)	9 (18%)	1 10
14	XN	49/50 (98%)	42 (86%)	7 (14%)	3 19
15	QO	79/80 (99%)	72 (91%)	7 (9%)	9 33
15	XO	79/80 (99%)	68 (86%)	11 (14%)	3 19
16	QP	72/74 (97%)	62 (86%)	10 (14%)	3 19
16	XP	72/74 (97%)	66 (92%)	6 (8%)	10 35
17	QQ	95/97 (98%)	86 (90%)	9 (10%)	8 32
17	XQ	95/97 (98%)	88 (93%)	7 (7%)	13 39
18	QR	61/77 (79%)	50 (82%)	11 (18%)	2 11
18	XR	61/77 (79%)	50 (82%)	11 (18%)	2 11
19	QS	73/80 (91%)	58 (80%)	15 (20%)	1 8
19	XS	73/80 (91%)	56 (77%)	17 (23%)	1 5
20	QT	76/82 (93%)	66 (87%)	10 (13%)	4 21
20	XT	76/82 (93%)	66 (87%)	10 (13%)	4 21
21	QU	20/22 (91%)	20 (100%)	0	100 100
21	XU	20/22 (91%)	20 (100%)	0	100 100
27	RD	214/218 (98%)	170 (79%)	44 (21%)	1 8
27	YD	214/218 (98%)	173 (81%)	41 (19%)	1 9
28	RE	165/166 (99%)	125 (76%)	40 (24%)	1 5
28	YE	165/166 (99%)	125 (76%)	40 (24%)	1 5
29	RF	161/166 (97%)	139 (86%)	22 (14%)	3 20
29	YF	161/166 (97%)	138 (86%)	23 (14%)	3 19
30	RG	155/156 (99%)	135 (87%)	20 (13%)	4 22
30	YG	155/156 (99%)	131 (84%)	24 (16%)	2 16
31	RH	142/148 (96%)	111 (78%)	31 (22%)	1 7
31	YH	142/148 (96%)	112 (79%)	30 (21%)	1 7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
32	RI	122/124 (98%)	100 (82%)	22 (18%)	2 11
32	YI	122/124 (98%)	92 (75%)	30 (25%)	1 4
33	RN	117/119 (98%)	94 (80%)	23 (20%)	1 8
33	YN	117/119 (98%)	94 (80%)	23 (20%)	1 8
34	RO	100/100 (100%)	91 (91%)	9 (9%)	9 33
34	YO	100/100 (100%)	85 (85%)	15 (15%)	3 17
35	RP	116/116 (100%)	85 (73%)	31 (27%)	0 3
35	YP	116/116 (100%)	78 (67%)	38 (33%)	0 2
36	RQ	111/111 (100%)	95 (86%)	16 (14%)	3 18
36	YQ	111/111 (100%)	94 (85%)	17 (15%)	3 16
37	RR	101/101 (100%)	81 (80%)	20 (20%)	1 8
37	YR	101/101 (100%)	80 (79%)	21 (21%)	1 7
38	RS	87/88 (99%)	67 (77%)	20 (23%)	1 6
38	YS	87/88 (99%)	73 (84%)	14 (16%)	2 15
39	RT	120/127 (94%)	101 (84%)	19 (16%)	2 15
39	YT	120/127 (94%)	94 (78%)	26 (22%)	1 7
40	RU	93/94 (99%)	79 (85%)	14 (15%)	3 17
40	YU	93/94 (99%)	79 (85%)	14 (15%)	3 17
41	RV	82/82 (100%)	63 (77%)	19 (23%)	1 5
41	YV	82/82 (100%)	66 (80%)	16 (20%)	1 9
42	RW	92/92 (100%)	73 (79%)	19 (21%)	1 7
42	YW	92/92 (100%)	76 (83%)	16 (17%)	2 12
43	RX	74/78 (95%)	61 (82%)	13 (18%)	2 12
43	YX	74/78 (95%)	59 (80%)	15 (20%)	1 8
44	RY	85/91 (93%)	56 (66%)	29 (34%)	0 1
44	YY	85/91 (93%)	62 (73%)	23 (27%)	0 3
45	RZ	162/179 (90%)	141 (87%)	21 (13%)	4 21
45	YZ	162/179 (90%)	141 (87%)	21 (13%)	4 21
46	R0	65/67 (97%)	58 (89%)	7 (11%)	6 28
46	Y0	65/67 (97%)	57 (88%)	8 (12%)	4 23
47	R1	82/83 (99%)	74 (90%)	8 (10%)	7 31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
47	Y1	82/83 (99%)	69 (84%)	13 (16%)	2 15
48	R2	64/67 (96%)	48 (75%)	16 (25%)	0 4
48	Y2	64/67 (96%)	57 (89%)	7 (11%)	6 27
49	R3	51/52 (98%)	43 (84%)	8 (16%)	2 15
49	Y3	51/52 (98%)	44 (86%)	7 (14%)	3 20
50	R4	63/63 (100%)	48 (76%)	15 (24%)	1 5
50	Y4	63/63 (100%)	43 (68%)	20 (32%)	0 2
51	R5	51/52 (98%)	40 (78%)	11 (22%)	1 7
51	Y5	51/52 (98%)	37 (72%)	14 (28%)	0 3
52	R6	48/52 (92%)	32 (67%)	16 (33%)	0 2
52	Y6	48/52 (92%)	34 (71%)	14 (29%)	0 2
53	R7	42/42 (100%)	35 (83%)	7 (17%)	2 14
53	Y7	42/42 (100%)	35 (83%)	7 (17%)	2 14
54	R8	54/55 (98%)	39 (72%)	15 (28%)	0 3
54	Y8	54/55 (98%)	39 (72%)	15 (28%)	0 3
55	R9	34/34 (100%)	30 (88%)	4 (12%)	5 24
55	Y9	34/34 (100%)	31 (91%)	3 (9%)	9 33
All	All	9702/10066 (96%)	8068 (83%)	1634 (17%)	2 13

All (1634) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	QB	5	ILE
2	QB	6	THR
2	QB	7	VAL
2	QB	8	LYS
2	QB	15	VAL
2	QB	23	ARG
2	QB	24	TRP
2	QB	32	ILE
2	QB	53	ARG
2	QB	60	ASP
2	QB	63	MET
2	QB	67	THR
2	QB	75	LYS
2	QB	76	GLN

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Mol	Chain	Res	Type
2	QB	87	ARG
2	QB	92	TYR
2	QB	94	ASN
2	QB	101	MET
2	QB	109	SER
2	QB	119	GLU
2	QB	121	LEU
2	QB	155	LEU
2	QB	158	LEU
2	QB	165	VAL
2	QB	168	THR
2	QB	170	GLU
2	QB	172	ILE
2	QB	175	ARG
2	QB	187	LEU
2	QB	196	LEU
2	QB	213	LEU
2	QB	215	LEU
2	QB	217	ARG
2	QB	221	LEU
2	QB	239	VAL
3	QC	3	ASN
3	QC	5	ILE
3	QC	12	LEU
3	QC	16	ARG
3	QC	21	ARG
3	QC	45	LYS
3	QC	76	VAL
3	QC	94	LEU
3	QC	154	SER
3	QC	162	GLN
3	QC	165	THR
3	QC	178	LEU
3	QC	206	GLU
4	QD	3	ARG
4	QD	9	CYS
4	QD	12	CYS
4	QD	14	ARG
4	QD	26	CYS
4	QD	30	LYS
4	QD	33	MET
4	QD	50	ARG

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Mol	Chain	Res	Type
4	QD	57	ARG
4	QD	58	LEU
4	QD	86	LYS
4	QD	96	LEU
4	QD	122	ARG
4	QD	127	THR
4	QD	131	ARG
4	QD	135	LEU
4	QD	154	ASN
4	QD	190	ASP
4	QD	192	GLU
5	QE	12	LEU
5	QE	13	ILE
5	QE	31	LEU
5	QE	34	VAL
5	QE	41	VAL
5	QE	51	VAL
5	QE	68	GLU
5	QE	71	LEU
5	QE	79	GLU
5	QE	81	GLU
5	QE	90	VAL
5	QE	98	THR
5	QE	101	ILE
5	QE	107	ARG
5	QE	115	VAL
5	QE	153	LYS
6	QF	16	GLN
6	QF	21	LEU
6	QF	37	VAL
6	QF	43	LEU
6	QF	45	LEU
6	QF	70	ASP
6	QF	72	VAL
6	QF	75	LEU
6	QF	81	ILE
6	QF	91	VAL
6	QF	98	LEU
7	QG	8	GLU
7	QG	38	LEU
7	QG	54	THR
7	QG	80	VAL

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Mol	Chain	Res	Type
7	QG	92	SER
7	QG	94	ARG
7	QG	104	LEU
7	QG	113	GLU
7	QG	135	VAL
7	QG	136	LYS
7	QG	137	LYS
8	QH	1	MET
8	QH	24	THR
8	QH	25	ASP
8	QH	26	VAL
8	QH	39	LEU
8	QH	41	ARG
8	QH	45	ILE
8	QH	54	ASP
8	QH	99	GLU
8	QH	109	ILE
8	QH	112	LEU
8	QH	121	ASP
8	QH	125	ARG
8	QH	129	VAL
9	QI	7	THR
9	QI	9	ARG
9	QI	10	ARG
9	QI	11	LYS
9	QI	23	ASN
9	QI	41	VAL
9	QI	47	LEU
9	QI	56	LEU
9	QI	64	THR
9	QI	65	VAL
9	QI	75	ASP
9	QI	86	VAL
9	QI	95	LYS
9	QI	104	ARG
9	QI	105	ASP
9	QI	114	TYR
9	QI	118	LYS
9	QI	128	ARG
10	QJ	3	LYS
10	QJ	16	LEU
10	QJ	22	LYS

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Mol	Chain	Res	Type
10	QJ	47	PHE
10	QJ	48	THR
10	QJ	57	LYS
10	QJ	62	HIS
10	QJ	74	ILE
10	QJ	80	LYS
10	QJ	84	GLN
10	QJ	92	THR
10	QJ	96	ILE
11	QK	21	ILE
11	QK	26	ASN
11	QK	29	ILE
11	QK	30	VAL
11	QK	32	ILE
11	QK	57	THR
11	QK	63	LEU
11	QK	92	GLU
11	QK	103	LEU
11	QK	109	VAL
11	QK	114	VAL
12	QL	17	LYS
12	QL	20	LYS
12	QL	24	VAL
12	QL	27	LEU
12	QL	41	ARG
12	QL	57	LYS
12	QL	62	SER
12	QL	70	ILE
12	QL	73	GLU
12	QL	81	SER
12	QL	101	VAL
12	QL	104	VAL
12	QL	110	VAL
12	QL	112	ASP
12	QL	116	SER
12	QL	120	TYR
12	QL	126	LYS
13	QM	8	GLU
13	QM	13	LYS
13	QM	17	VAL
13	QM	19	LEU
13	QM	45	VAL

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Mol	Chain	Res	Type
13	QM	47	ASP
13	QM	48	LEU
13	QM	56	LEU
13	QM	57	ARG
13	QM	64	TRP
13	QM	66	LEU
13	QM	70	LEU
13	QM	77	ASN
13	QM	78	ILE
13	QM	84	ILE
13	QM	88	ARG
13	QM	90	LEU
13	QM	98	VAL
13	QM	108	ARG
13	QM	111	LYS
13	QM	114	ARG
13	QM	115	LYS
13	QM	117	VAL
13	QM	122	LYS
14	QN	6	LEU
14	QN	12	ARG
14	QN	13	THR
14	QN	18	VAL
14	QN	33	VAL
14	QN	43	CYS
14	QN	44	LEU
14	QN	46	GLU
14	QN	57	ARG
15	QO	3	ILE
15	QO	4	THR
15	QO	26	GLU
15	QO	31	LEU
15	QO	39	LEU
15	QO	64	ARG
15	QO	82	ILE
16	QP	2	VAL
16	QP	20	VAL
16	QP	26	ARG
16	QP	28	ARG
16	QP	33	ILE
16	QP	53	VAL
16	QP	67	THR

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Mol	Chain	Res	Type
16	QP	69	THR
16	QP	71	ARG
16	QP	74	LEU
17	QQ	37	LYS
17	QQ	38	ARG
17	QQ	43	LEU
17	QQ	52	LYS
17	QQ	59	ILE
17	QQ	62	SER
17	QQ	68	ARG
17	QQ	74	LEU
17	QQ	101	ARG
18	QR	26	LEU
18	QR	29	PHE
18	QR	31	LEU
18	QR	32	ARG
18	QR	36	ASN
18	QR	46	GLU
18	QR	54	ARG
18	QR	76	LEU
18	QR	82	THR
18	QR	83	GLU
18	QR	86	VAL
19	QS	5	LEU
19	QS	10	PHE
19	QS	11	VAL
19	QS	12	ASP
19	QS	21	GLU
19	QS	28	LYS
19	QS	29	ARG
19	QS	30	LEU
19	QS	37	ARG
19	QS	43	GLU
19	QS	62	ILE
19	QS	63	THR
19	QS	67	VAL
19	QS	77	THR
19	QS	83	HIS
20	QT	17	ARG
20	QT	24	LEU
20	QT	41	ILE
20	QT	45	GLN

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Mol	Chain	Res	Type
20	QT	72	LEU
20	QT	73	HIS
20	QT	75	ASN
20	QT	84	LEU
20	QT	93	GLU
20	QT	99	LEU
27	RD	3	VAL
27	RD	5	LYS
27	RD	10	THR
27	RD	17	THR
27	RD	25	THR
27	RD	34	VAL
27	RD	40	THR
27	RD	44	ASN
27	RD	46	GLN
27	RD	49	ILE
27	RD	61	LEU
27	RD	64	ILE
27	RD	65	ILE
27	RD	69	ARG
27	RD	71	ASP
27	RD	73	VAL
27	RD	83	GLU
27	RD	87	ASN
27	RD	89	SER
27	RD	95	LEU
27	RD	98	VAL
27	RD	103	ARG
27	RD	105	ILE
27	RD	106	ILE
27	RD	111	LEU
27	RD	134	ARG
27	RD	136	ILE
27	RD	147	LEU
27	RD	150	LYS
27	RD	155	LEU
27	RD	157	ARG
27	RD	169	GLU
27	RD	173	VAL
27	RD	192	THR
27	RD	193	VAL
27	RD	212	SER

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Mol	Chain	Res	Type
27	RD	221	VAL
27	RD	229	VAL
27	RD	237	GLU
27	RD	257	LEU
27	RD	259	THR
27	RD	261	LYS
27	RD	271	ILE
27	RD	273	ARG
28	RE	2	LYS
28	RE	4	ILE
28	RE	5	LEU
28	RE	13	ARG
28	RE	14	ILE
28	RE	17	ASP
28	RE	21	VAL
28	RE	25	VAL
28	RE	26	ILE
28	RE	27	LEU
28	RE	33	VAL
28	RE	36	ARG
28	RE	38	THR
28	RE	41	LYS
28	RE	45	THR
28	RE	48	GLN
28	RE	49	LEU
28	RE	54	GLN
28	RE	59	VAL
28	RE	61	ARG
28	RE	62	PRO
28	RE	66	HIS
28	RE	73	GLU
28	RE	75	VAL
28	RE	77	ILE
28	RE	78	LEU
28	RE	79	ARG
28	RE	80	GLU
28	RE	93	VAL
28	RE	117	MET
28	RE	119	ARG
28	RE	143	ASN
28	RE	146	THR
28	RE	154	LYS

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Mol	Chain	Res	Type
28	RE	167	VAL
28	RE	179	GLU
28	RE	184	VAL
28	RE	196	VAL
28	RE	200	GLU
28	RE	203	LYS
29	RF	9	ILE
29	RF	11	VAL
29	RF	13	SER
29	RF	28	ILE
29	RF	32	LEU
29	RF	33	LEU
29	RF	45	ARG
29	RF	57	VAL
29	RF	65	TRP
29	RF	70	THR
29	RF	74	ARG
29	RF	78	ILE
29	RF	84	VAL
29	RF	108	LYS
29	RF	117	ARG
29	RF	174	VAL
29	RF	181	LEU
29	RF	183	VAL
29	RF	192	LEU
29	RF	194	MET
29	RF	197	ASP
29	RF	206	ILE
30	RG	7	LEU
30	RG	10	LYS
30	RG	20	ILE
30	RG	26	GLN
30	RG	34	LEU
30	RG	43	LEU
30	RG	45	GLU
30	RG	53	LEU
30	RG	54	GLU
30	RG	63	ILE
30	RG	67	LYS
30	RG	71	THR
30	RG	88	ILE
30	RG	94	LEU

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Mol	Chain	Res	Type
30	RG	98	ARG
30	RG	118	ARG
30	RG	147	ASP
30	RG	159	VAL
30	RG	167	GLU
30	RG	174	GLU
31	RH	3	ARG
31	RH	4	ILE
31	RH	9	ILE
31	RH	10	PRO
31	RH	11	VAL
31	RH	16	SER
31	RH	27	LYS
31	RH	32	GLU
31	RH	37	VAL
31	RH	41	MET
31	RH	43	VAL
31	RH	44	VAL
31	RH	47	GLU
31	RH	59	ARG
31	RH	64	LEU
31	RH	77	LYS
31	RH	81	GLU
31	RH	88	LEU
31	RH	89	ILE
31	RH	105	LEU
31	RH	124	GLU
31	RH	132	ARG
31	RH	136	ILE
31	RH	139	GLN
31	RH	143	GLN
31	RH	144	VAL
31	RH	152	ARG
31	RH	153	LYS
31	RH	154	PRO
31	RH	155	SER
31	RH	169	VAL
32	RI	2	LYS
32	RI	7	GLU
32	RI	9	LEU
32	RI	10	GLU
32	RI	27	ARG

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Mol	Chain	Res	Type
32	RI	33	ARG
32	RI	38	LEU
32	RI	44	LEU
32	RI	56	LYS
32	RI	57	ARG
32	RI	81	VAL
32	RI	85	GLU
32	RI	86	THR
32	RI	92	VAL
32	RI	97	ILE
32	RI	101	LEU
32	RI	128	LEU
32	RI	129	THR
32	RI	134	PRO
32	RI	135	GLU
32	RI	139	GLN
32	RI	142	VAL
33	RN	1	MET
33	RN	2	LYS
33	RN	5	VAL
33	RN	7	LYS
33	RN	12	ARG
33	RN	14	VAL
33	RN	30	ILE
33	RN	32	THR
33	RN	33	LEU
33	RN	34	LEU
33	RN	43	THR
33	RN	48	MET
33	RN	60	ILE
33	RN	61	ARG
33	RN	62	VAL
33	RN	87	LEU
33	RN	90	MET
33	RN	96	GLU
33	RN	98	VAL
33	RN	107	LEU
33	RN	109	LYS
33	RN	120	LEU
33	RN	136	GLU
34	RO	9	GLU
34	RO	19	ILE

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Mol	Chain	Res	Type
34	RO	24	VAL
34	RO	63	VAL
34	RO	69	ILE
34	RO	91	LEU
34	RO	98	VAL
34	RO	102	VAL
34	RO	115	VAL
35	RP	6	LEU
35	RP	7	ARG
35	RP	14	LYS
35	RP	15	ARG
35	RP	18	ARG
35	RP	19	VAL
35	RP	21	ARG
35	RP	30	THR
35	RP	36	LYS
35	RP	41	ARG
35	RP	45	LEU
35	RP	50	ARG
35	RP	59	LEU
35	RP	62	LEU
35	RP	67	MET
35	RP	68	GLN
35	RP	71	VAL
35	RP	75	ILE
35	RP	81	GLN
35	RP	88	LEU
35	RP	91	PHE
35	RP	100	LEU
35	RP	105	LEU
35	RP	106	LEU
35	RP	111	ARG
35	RP	112	LEU
35	RP	135	LEU
35	RP	138	LEU
35	RP	139	LYS
35	RP	144	GLU
35	RP	146	VAL
36	RQ	2	LEU
36	RQ	25	ASP
36	RQ	26	TYR
36	RQ	27	VAL

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Mol	Chain	Res	Type
36	RQ	46	GLN
36	RQ	54	MET
36	RQ	55	VAL
36	RQ	58	PHE
36	RQ	60	ARG
36	RQ	79	LEU
36	RQ	83	MET
36	RQ	89	ASN
36	RQ	90	VAL
36	RQ	91	GLU
36	RQ	112	GLU
36	RQ	139	GLU
37	RR	1	MET
37	RR	6	SER
37	RR	9	LYS
37	RR	18	LEU
37	RR	28	LEU
37	RR	29	LEU
37	RR	35	THR
37	RR	44	LEU
37	RR	45	ARG
37	RR	67	LEU
37	RR	71	GLN
37	RR	73	VAL
37	RR	75	LEU
37	RR	79	LEU
37	RR	91	GLN
37	RR	95	THR
37	RR	100	LEU
37	RR	104	ARG
37	RR	117	VAL
37	RR	118	GLU
38	RS	4	LEU
38	RS	12	PHE
38	RS	17	ARG
38	RS	18	ILE
38	RS	20	ARG
38	RS	24	LEU
38	RS	27	SER
38	RS	39	ILE
38	RS	44	LYS
38	RS	50	SER

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Mol	Chain	Res	Type
38	RS	54	LEU
38	RS	56	LEU
38	RS	57	LYS
38	RS	58	LEU
38	RS	59	LYS
38	RS	69	VAL
38	RS	98	VAL
38	RS	101	LEU
38	RS	103	GLU
38	RS	106	ARG
39	RT	7	ILE
39	RT	18	ASP
39	RT	27	THR
39	RT	30	VAL
39	RT	41	ARG
39	RT	42	ILE
39	RT	50	ILE
39	RT	51	ARG
39	RT	62	THR
39	RT	63	VAL
39	RT	65	LYS
39	RT	74	ARG
39	RT	78	LEU
39	RT	88	ILE
39	RT	89	VAL
39	RT	99	LEU
39	RT	112	ARG
39	RT	125	ARG
39	RT	128	GLU
40	RU	55	ARG
40	RU	59	ARG
40	RU	60	LEU
40	RU	64	ARG
40	RU	65	ILE
40	RU	74	LEU
40	RU	90	VAL
40	RU	92	ARG
40	RU	98	LEU
40	RU	100	VAL
40	RU	108	GLU
40	RU	111	GLU
40	RU	114	LYS

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Mol	Chain	Res	Type
40	RU	117	GLN
41	RV	19	LYS
41	RV	21	ARG
41	RV	22	VAL
41	RV	24	LYS
41	RV	35	LEU
41	RV	37	VAL
41	RV	38	LEU
41	RV	45	THR
41	RV	46	VAL
41	RV	47	VAL
41	RV	49	THR
41	RV	57	VAL
41	RV	61	VAL
41	RV	62	LEU
41	RV	72	VAL
41	RV	74	LYS
41	RV	78	LYS
41	RV	79	VAL
41	RV	99	ILE
42	RW	11	ARG
42	RW	16	LYS
42	RW	19	LEU
42	RW	20	VAL
42	RW	21	VAL
42	RW	23	LEU
42	RW	27	LYS
42	RW	30	GLU
42	RW	51	LEU
42	RW	60	ASN
42	RW	63	ASP
42	RW	67	ASP
42	RW	76	VAL
42	RW	82	LEU
42	RW	92	ARG
42	RW	95	ILE
42	RW	100	THR
42	RW	106	ILE
42	RW	107	LEU
43	RX	12	VAL
43	RX	15	GLU
43	RX	23	GLU

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Mol	Chain	Res	Type
43	RX	27	THR
43	RX	30	VAL
43	RX	35	THR
43	RX	49	VAL
43	RX	52	VAL
43	RX	54	VAL
43	RX	65	ARG
43	RX	73	ARG
43	RX	80	ILE
43	RX	81	VAL
44	RY	3	VAL
44	RY	13	VAL
44	RY	14	LEU
44	RY	26	LYS
44	RY	27	VAL
44	RY	30	VAL
44	RY	34	LYS
44	RY	37	VAL
44	RY	38	ILE
44	RY	39	VAL
44	RY	43	ASN
44	RY	45	VAL
44	RY	51	VAL
44	RY	55	TYR
44	RY	57	GLN
44	RY	61	ILE
44	RY	64	GLU
44	RY	67	LEU
44	RY	70	SER
44	RY	71	LYS
44	RY	75	ILE
44	RY	76	CYS
44	RY	87	LYS
44	RY	88	LYS
44	RY	90	LEU
44	RY	95	LYS
44	RY	96	ILE
44	RY	97	ARG
44	RY	102	CYS
45	RZ	2	GLU
45	RZ	5	LEU
45	RZ	19	ARG

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Mol	Chain	Res	Type
45	RZ	20	ARG
45	RZ	52	SER
45	RZ	60	GLU
45	RZ	70	LEU
45	RZ	87	ASP
45	RZ	94	GLU
45	RZ	111	VAL
45	RZ	121	HIS
45	RZ	128	VAL
45	RZ	145	GLU
45	RZ	150	LEU
45	RZ	151	HIS
45	RZ	163	LEU
45	RZ	166	SER
45	RZ	168	GLU
45	RZ	174	VAL
45	RZ	182	LYS
45	RZ	183	LEU
46	R0	5	LYS
46	R0	7	LEU
46	R0	10	THR
46	R0	36	ILE
46	R0	53	MET
46	R0	66	VAL
46	R0	74	ARG
47	R1	7	ILE
47	R1	21	ARG
47	R1	41	ARG
47	R1	51	VAL
47	R1	62	VAL
47	R1	80	LEU
47	R1	90	ILE
47	R1	91	LYS
48	R2	5	GLU
48	R2	17	SER
48	R2	22	GLU
48	R2	23	LYS
48	R2	24	LEU
48	R2	27	GLU
48	R2	32	LEU
48	R2	41	ILE
48	R2	45	SER

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Mol	Chain	Res	Type
48	R2	47	ASN
48	R2	48	HIS
48	R2	53	LEU
48	R2	62	THR
48	R2	63	VAL
48	R2	65	ASN
48	R2	71	ASN
49	R3	6	VAL
49	R3	8	LEU
49	R3	18	ASP
49	R3	32	GLN
49	R3	40	THR
49	R3	48	GLU
49	R3	56	VAL
49	R3	59	VAL
50	R4	6	HIS
50	R4	15	ILE
50	R4	21	VAL
50	R4	23	GLU
50	R4	42	PHE
50	R4	49	PHE
50	R4	50	VAL
50	R4	53	GLU
50	R4	57	GLU
50	R4	61	ARG
50	R4	62	ARG
50	R4	63	TYR
50	R4	68	ARG
50	R4	69	LYS
50	R4	71	ARG
51	R5	3	LYS
51	R5	4	HIS
51	R5	6	VAL
51	R5	11	THR
51	R5	19	ARG
51	R5	25	LEU
51	R5	36	CYS
51	R5	37	LYS
51	R5	52	TYR
51	R5	56	LYS
51	R5	58	LEU
52	R6	6	ARG

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Mol	Chain	Res	Type
52	R6	8	LYS
52	R6	9	LEU
52	R6	10	LEU
52	R6	11	LEU
52	R6	17	LYS
52	R6	19	ARG
52	R6	23	THR
52	R6	27	LYS
52	R6	30	THR
52	R6	34	LEU
52	R6	35	GLU
52	R6	37	ARG
52	R6	44	ARG
52	R6	47	THR
52	R6	53	LYS
53	R7	1	MET
53	R7	2	LYS
53	R7	4	THR
53	R7	9	ARG
53	R7	14	LYS
53	R7	43	THR
53	R7	46	VAL
54	R8	3	LYS
54	R8	16	ILE
54	R8	30	ARG
54	R8	35	GLN
54	R8	39	LYS
54	R8	43	GLN
54	R8	47	LYS
54	R8	48	PHE
54	R8	49	VAL
54	R8	52	LYS
54	R8	53	PRO
54	R8	59	LYS
54	R8	62	LEU
54	R8	63	PRO
54	R8	65	GLU
55	R9	1	MET
55	R9	17	ILE
55	R9	29	ASN
55	R9	33	LYS
2	XB	5	ILE

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Mol	Chain	Res	Type
2	XB	7	VAL
2	XB	8	LYS
2	XB	15	VAL
2	XB	23	ARG
2	XB	36	ARG
2	XB	67	THR
2	XB	68	ILE
2	XB	71	VAL
2	XB	74	LYS
2	XB	75	LYS
2	XB	92	TYR
2	XB	136	VAL
2	XB	145	LEU
2	XB	155	LEU
2	XB	172	ILE
2	XB	175	ARG
2	XB	178	ARG
2	XB	187	LEU
2	XB	195	ASP
2	XB	196	LEU
2	XB	200	ILE
2	XB	215	LEU
2	XB	223	ILE
2	XB	235	SER
3	XC	3	ASN
3	XC	5	ILE
3	XC	12	LEU
3	XC	14	ILE
3	XC	15	THR
3	XC	21	ARG
3	XC	27	LYS
3	XC	47	LEU
3	XC	56	ASP
3	XC	76	VAL
3	XC	95	THR
3	XC	162	GLN
3	XC	178	LEU
3	XC	192	THR
4	XD	3	ARG
4	XD	9	CYS
4	XD	15	GLU
4	XD	19	LEU

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Mol	Chain	Res	Type
4	XD	30	LYS
4	XD	33	MET
4	XD	50	ARG
4	XD	53	ASP
4	XD	58	LEU
4	XD	76	ARG
4	XD	84	LYS
4	XD	86	LYS
4	XD	88	VAL
4	XD	92	VAL
4	XD	96	LEU
4	XD	108	LEU
4	XD	122	ARG
4	XD	127	THR
4	XD	131	ARG
4	XD	137	SER
4	XD	150	GLU
4	XD	154	ASN
4	XD	175	SER
4	XD	190	ASP
4	XD	208	SER
5	XE	6	PHE
5	XE	7	GLU
5	XE	8	GLU
5	XE	10	MET
5	XE	11	ILE
5	XE	13	ILE
5	XE	18	ARG
5	XE	31	LEU
5	XE	41	VAL
5	XE	73	ASN
5	XE	76	ILE
5	XE	79	GLU
5	XE	89	ILE
5	XE	101	ILE
5	XE	147	ASP
5	XE	153	LYS
6	XF	10	LEU
6	XF	21	LEU
6	XF	23	LYS
6	XF	36	ARG
6	XF	37	VAL

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Mol	Chain	Res	Type
6	XF	45	LEU
6	XF	71	ARG
6	XF	72	VAL
6	XF	74	ASP
6	XF	91	VAL
6	XF	92	LYS
6	XF	98	LEU
7	XG	5	ARG
7	XG	8	GLU
7	XG	13	GLN
7	XG	27	ILE
7	XG	35	LYS
7	XG	54	THR
7	XG	63	LYS
7	XG	78	ARG
7	XG	80	VAL
7	XG	104	LEU
7	XG	113	GLU
7	XG	114	ARG
7	XG	136	LYS
7	XG	137	LYS
8	XH	1	MET
8	XH	12	ARG
8	XH	19	VAL
8	XH	24	THR
8	XH	26	VAL
8	XH	41	ARG
8	XH	54	ASP
8	XH	63	LEU
8	XH	80	ILE
8	XH	109	ILE
8	XH	112	LEU
8	XH	121	ASP
8	XH	137	VAL
9	XI	9	ARG
9	XI	41	VAL
9	XI	44	VAL
9	XI	56	LEU
9	XI	63	ILE
9	XI	65	VAL
9	XI	70	LYS
9	XI	79	LEU

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Mol	Chain	Res	Type
9	XI	95	LYS
9	XI	96	LEU
9	XI	99	LEU
9	XI	102	LEU
9	XI	104	ARG
9	XI	105	ASP
9	XI	108	VAL
9	XI	111	ARG
9	XI	112	LYS
9	XI	114	TYR
9	XI	118	LYS
9	XI	124	GLN
9	XI	128	ARG
10	XJ	3	LYS
10	XJ	17	ASP
10	XJ	45	ARG
10	XJ	47	PHE
10	XJ	49	VAL
10	XJ	57	LYS
10	XJ	62	HIS
10	XJ	70	ARG
10	XJ	72	VAL
10	XJ	74	ILE
10	XJ	80	LYS
10	XJ	84	GLN
10	XJ	90	LEU
10	XJ	96	ILE
10	XJ	98	ILE
11	XK	26	ASN
11	XK	29	ILE
11	XK	31	THR
11	XK	32	ILE
11	XK	36	ASP
11	XK	57	THR
11	XK	96	ARG
11	XK	114	VAL
12	XL	17	LYS
12	XL	20	LYS
12	XL	24	VAL
12	XL	27	LEU
12	XL	41	ARG
12	XL	48	PRO

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Mol	Chain	Res	Type
12	XL	57	LYS
12	XL	62	SER
12	XL	70	ILE
12	XL	73	GLU
12	XL	81	SER
12	XL	101	VAL
12	XL	104	VAL
12	XL	110	VAL
12	XL	112	ASP
12	XL	116	SER
12	XL	120	TYR
12	XL	126	LYS
13	XM	3	ARG
13	XM	13	LYS
13	XM	17	VAL
13	XM	19	LEU
13	XM	32	GLU
13	XM	45	VAL
13	XM	48	LEU
13	XM	56	LEU
13	XM	64	TRP
13	XM	66	LEU
13	XM	70	LEU
13	XM	84	ILE
13	XM	88	ARG
13	XM	98	VAL
13	XM	114	ARG
13	XM	115	LYS
13	XM	117	VAL
13	XM	122	LYS
14	XN	6	LEU
14	XN	12	ARG
14	XN	32	SER
14	XN	33	VAL
14	XN	40	CYS
14	XN	41	ARG
14	XN	44	LEU
15	XO	3	ILE
15	XO	6	GLU
15	XO	8	LYS
15	XO	24	SER
15	XO	26	GLU

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Mol	Chain	Res	Type
15	XO	39	LEU
15	XO	47	LYS
15	XO	65	ARG
15	XO	66	LEU
15	XO	82	ILE
15	XO	87	ILE
16	XP	2	VAL
16	XP	11	SER
16	XP	20	VAL
16	XP	28	ARG
16	XP	67	THR
16	XP	69	THR
17	XQ	9	VAL
17	XQ	52	LYS
17	XQ	59	ILE
17	XQ	62	SER
17	XQ	68	ARG
17	XQ	74	LEU
17	XQ	101	ARG
18	XR	26	LEU
18	XR	29	PHE
18	XR	32	ARG
18	XR	36	ASN
18	XR	41	LYS
18	XR	46	GLU
18	XR	54	ARG
18	XR	76	LEU
18	XR	82	THR
18	XR	83	GLU
18	XR	86	VAL
19	XS	5	LEU
19	XS	10	PHE
19	XS	11	VAL
19	XS	12	ASP
19	XS	13	ASP
19	XS	21	GLU
19	XS	22	LEU
19	XS	29	ARG
19	XS	30	LEU
19	XS	31	ILE
19	XS	37	ARG
19	XS	44	MET

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Mol	Chain	Res	Type
19	XS	58	VAL
19	XS	63	THR
19	XS	78	ARG
19	XS	81	ARG
19	XS	83	HIS
20	XT	10	LEU
20	XT	13	LEU
20	XT	24	LEU
20	XT	41	ILE
20	XT	45	GLN
20	XT	50	GLU
20	XT	55	ILE
20	XT	73	HIS
20	XT	84	LEU
20	XT	93	GLU
27	YD	10	THR
27	YD	17	THR
27	YD	26	LYS
27	YD	28	GLU
27	YD	33	LEU
27	YD	43	ARG
27	YD	44	ASN
27	YD	52	ARG
27	YD	61	LEU
27	YD	65	ILE
27	YD	67	PHE
27	YD	73	VAL
27	YD	94	LEU
27	YD	98	VAL
27	YD	105	ILE
27	YD	106	ILE
27	YD	131	LEU
27	YD	134	ARG
27	YD	135	PHE
27	YD	147	LEU
27	YD	155	LEU
27	YD	157	ARG
27	YD	166	GLN
27	YD	173	VAL
27	YD	177	LEU
27	YD	183	ARG
27	YD	192	THR

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Mol	Chain	Res	Type
27	YD	198	ASN
27	YD	200	ASP
27	YD	215	LEU
27	YD	217	ARG
27	YD	218	ARG
27	YD	221	VAL
27	YD	226	MET
27	YD	230	ASP
27	YD	233	HIS
27	YD	237	GLU
27	YD	257	LEU
27	YD	259	THR
27	YD	261	LYS
27	YD	271	ILE
28	YE	2	LYS
28	YE	4	ILE
28	YE	5	LEU
28	YE	13	ARG
28	YE	14	ILE
28	YE	17	ASP
28	YE	21	VAL
28	YE	25	VAL
28	YE	26	ILE
28	YE	27	LEU
28	YE	33	VAL
28	YE	36	ARG
28	YE	38	THR
28	YE	41	LYS
28	YE	45	THR
28	YE	48	GLN
28	YE	49	LEU
28	YE	54	GLN
28	YE	59	VAL
28	YE	61	ARG
28	YE	62	PRO
28	YE	66	HIS
28	YE	73	GLU
28	YE	75	VAL
28	YE	77	ILE
28	YE	78	LEU
28	YE	79	ARG
28	YE	80	GLU

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Mol	Chain	Res	Type
28	YE	93	VAL
28	YE	117	MET
28	YE	119	ARG
28	YE	143	ASN
28	YE	146	THR
28	YE	154	LYS
28	YE	167	VAL
28	YE	179	GLU
28	YE	184	VAL
28	YE	196	VAL
28	YE	200	GLU
28	YE	203	LYS
29	YF	7	TYR
29	YF	9	ILE
29	YF	25	PRO
29	YF	32	LEU
29	YF	33	LEU
29	YF	45	ARG
29	YF	46	ARG
29	YF	65	TRP
29	YF	66	PRO
29	YF	67	GLN
29	YF	70	THR
29	YF	82	ILE
29	YF	106	ARG
29	YF	108	LYS
29	YF	124	LEU
29	YF	127	GLU
29	YF	145	GLU
29	YF	161	GLU
29	YF	164	ARG
29	YF	179	GLU
29	YF	181	LEU
29	YF	183	VAL
29	YF	206	ILE
30	YG	3	LEU
30	YG	7	LEU
30	YG	19	LEU
30	YG	22	ARG
30	YG	31	VAL
30	YG	34	LEU
30	YG	37	VAL

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Mol	Chain	Res	Type
30	YG	43	LEU
30	YG	45	GLU
30	YG	58	GLN
30	YG	63	ILE
30	YG	64	THR
30	YG	66	GLN
30	YG	67	LYS
30	YG	80	PHE
30	YG	82	LEU
30	YG	84	LYS
30	YG	88	ILE
30	YG	90	LEU
30	YG	94	LEU
30	YG	118	ARG
30	YG	145	THR
30	YG	159	VAL
30	YG	167	GLU
31	YH	3	ARG
31	YH	4	ILE
31	YH	9	ILE
31	YH	10	PRO
31	YH	11	VAL
31	YH	16	SER
31	YH	27	LYS
31	YH	32	GLU
31	YH	37	VAL
31	YH	41	MET
31	YH	43	VAL
31	YH	44	VAL
31	YH	59	ARG
31	YH	64	LEU
31	YH	77	LYS
31	YH	81	GLU
31	YH	88	LEU
31	YH	89	ILE
31	YH	105	LEU
31	YH	124	GLU
31	YH	132	ARG
31	YH	136	ILE
31	YH	139	GLN
31	YH	143	GLN
31	YH	144	VAL

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Mol	Chain	Res	Type
31	YH	152	ARG
31	YH	153	LYS
31	YH	154	PRO
31	YH	155	SER
31	YH	169	VAL
32	YI	1	MET
32	YI	2	LYS
32	YI	6	LEU
32	YI	10	GLU
32	YI	20	ASP
32	YI	35	LEU
32	YI	37	VAL
32	YI	38	LEU
32	YI	40	THR
32	YI	56	LYS
32	YI	67	ARG
32	YI	70	GLU
32	YI	72	LEU
32	YI	77	LEU
32	YI	81	VAL
32	YI	85	GLU
32	YI	86	THR
32	YI	92	VAL
32	YI	95	LYS
32	YI	101	LEU
32	YI	112	LYS
32	YI	113	ARG
32	YI	118	LYS
32	YI	131	LYS
32	YI	135	GLU
32	YI	136	VAL
32	YI	138	ILE
32	YI	139	GLN
32	YI	142	VAL
32	YI	145	VAL
33	YN	2	LYS
33	YN	5	VAL
33	YN	7	LYS
33	YN	9	VAL
33	YN	15	LEU
33	YN	32	THR
33	YN	34	LEU

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Mol	Chain	Res	Type
33	YN	43	THR
33	YN	60	ILE
33	YN	61	ARG
33	YN	62	VAL
33	YN	65	LYS
33	YN	67	LEU
33	YN	73	THR
33	YN	90	MET
33	YN	96	GLU
33	YN	98	VAL
33	YN	99	LEU
33	YN	109	LYS
33	YN	112	LEU
33	YN	116	LEU
33	YN	120	LEU
33	YN	136	GLU
34	YO	2	ILE
34	YO	9	GLU
34	YO	19	ILE
34	YO	20	MET
34	YO	23	ARG
34	YO	24	VAL
34	YO	47	ILE
34	YO	49	ARG
34	YO	53	LYS
34	YO	66	LYS
34	YO	68	GLU
34	YO	85	VAL
34	YO	91	LEU
34	YO	98	VAL
34	YO	106	LEU
35	YP	4	SER
35	YP	6	LEU
35	YP	7	ARG
35	YP	13	ASN
35	YP	14	LYS
35	YP	15	ARG
35	YP	16	ARG
35	YP	19	VAL
35	YP	21	ARG
35	YP	27	HIS
35	YP	29	LYS

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Mol	Chain	Res	Type
35	YP	32	THR
35	YP	36	LYS
35	YP	45	LEU
35	YP	50	ARG
35	YP	61	ARG
35	YP	65	ARG
35	YP	71	VAL
35	YP	75	ILE
35	YP	81	GLN
35	YP	88	LEU
35	YP	91	PHE
35	YP	98	GLU
35	YP	100	LEU
35	YP	101	VAL
35	YP	105	LEU
35	YP	106	LEU
35	YP	112	LEU
35	YP	115	LEU
35	YP	123	LEU
35	YP	125	VAL
35	YP	126	VAL
35	YP	135	LEU
35	YP	139	LYS
35	YP	144	GLU
35	YP	146	VAL
35	YP	147	LEU
35	YP	149	GLU
36	YQ	2	LEU
36	YQ	14	ARG
36	YQ	25	ASP
36	YQ	26	TYR
36	YQ	27	VAL
36	YQ	46	GLN
36	YQ	54	MET
36	YQ	55	VAL
36	YQ	59	ARG
36	YQ	79	LEU
36	YQ	83	MET
36	YQ	89	ASN
36	YQ	90	VAL
36	YQ	91	GLU
36	YQ	106	VAL

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Mol	Chain	Res	Type
36	YQ	112	GLU
36	YQ	139	GLU
37	YR	18	LEU
37	YR	28	LEU
37	YR	29	LEU
37	YR	34	ILE
37	YR	36	THR
37	YR	40	LYS
37	YR	44	LEU
37	YR	51	LEU
37	YR	54	LEU
37	YR	57	ARG
37	YR	65	LEU
37	YR	67	LEU
37	YR	75	LEU
37	YR	79	LEU
37	YR	83	ILE
37	YR	95	THR
37	YR	99	LYS
37	YR	100	LEU
37	YR	102	GLU
37	YR	104	ARG
37	YR	105	ARG
38	YS	4	LEU
38	YS	12	PHE
38	YS	17	ARG
38	YS	18	ILE
38	YS	20	ARG
38	YS	44	LYS
38	YS	56	LEU
38	YS	57	LYS
38	YS	73	LEU
38	YS	89	ARG
38	YS	101	LEU
38	YS	103	GLU
38	YS	106	ARG
38	YS	111	GLU
39	YT	15	VAL
39	YT	17	THR
39	YT	23	ARG
39	YT	27	THR
39	YT	28	VAL

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Mol	Chain	Res	Type
39	YT	40	THR
39	YT	41	ARG
39	YT	42	ILE
39	YT	50	ILE
39	YT	65	LYS
39	YT	66	VAL
39	YT	74	ARG
39	YT	78	LEU
39	YT	85	LYS
39	YT	86	ILE
39	YT	87	ASP
39	YT	88	ILE
39	YT	89	VAL
39	YT	99	LEU
39	YT	110	ILE
39	YT	112	ARG
39	YT	115	ARG
39	YT	121	ILE
39	YT	125	ARG
39	YT	128	GLU
39	YT	134	GLU
40	YU	5	LYS
40	YU	11	ARG
40	YU	27	LEU
40	YU	51	LYS
40	YU	52	ARG
40	YU	60	LEU
40	YU	64	ARG
40	YU	74	LEU
40	YU	88	ILE
40	YU	90	VAL
40	YU	98	LEU
40	YU	104	GLN
40	YU	108	GLU
40	YU	111	GLU
41	YV	6	LYS
41	YV	7	THR
41	YV	10	LYS
41	YV	19	LYS
41	YV	32	THR
41	YV	35	LEU
41	YV	39	LEU

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Mol	Chain	Res	Type
41	YV	45	THR
41	YV	47	VAL
41	YV	57	VAL
41	YV	61	VAL
41	YV	66	ARG
41	YV	72	VAL
41	YV	78	LYS
41	YV	79	VAL
41	YV	99	ILE
42	YW	11	ARG
42	YW	16	LYS
42	YW	19	LEU
42	YW	23	LEU
42	YW	37	ARG
42	YW	51	LEU
42	YW	67	ASP
42	YW	69	LEU
42	YW	76	VAL
42	YW	88	ARG
42	YW	92	ARG
42	YW	95	ILE
42	YW	96	ILE
42	YW	100	THR
42	YW	106	ILE
42	YW	107	LEU
43	YX	12	VAL
43	YX	15	GLU
43	YX	27	THR
43	YX	36	LYS
43	YX	43	VAL
43	YX	49	VAL
43	YX	52	VAL
43	YX	57	LEU
43	YX	59	VAL
43	YX	63	LYS
43	YX	65	ARG
43	YX	66	LEU
43	YX	80	ILE
43	YX	83	VAL
43	YX	88	LYS
44	YY	14	LEU
44	YY	26	LYS

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Mol	Chain	Res	Type
44	YY	27	VAL
44	YY	28	LYS
44	YY	29	GLU
44	YY	34	LYS
44	YY	38	ILE
44	YY	44	ILE
44	YY	45	VAL
44	YY	51	VAL
44	YY	52	SER
44	YY	57	GLN
44	YY	61	ILE
44	YY	64	GLU
44	YY	67	LEU
44	YY	71	LYS
44	YY	73	ARG
44	YY	75	ILE
44	YY	84	ARG
44	YY	87	LYS
44	YY	90	LEU
44	YY	95	LYS
44	YY	97	ARG
45	YZ	2	GLU
45	YZ	5	LEU
45	YZ	19	ARG
45	YZ	20	ARG
45	YZ	28	MET
45	YZ	41	LEU
45	YZ	42	VAL
45	YZ	53	ILE
45	YZ	60	GLU
45	YZ	70	LEU
45	YZ	71	VAL
45	YZ	76	LEU
45	YZ	87	ASP
45	YZ	94	GLU
45	YZ	111	VAL
45	YZ	124	ILE
45	YZ	140	ASP
45	YZ	141	VAL
45	YZ	144	LEU
45	YZ	150	LEU
45	YZ	151	HIS

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Mol	Chain	Res	Type
46	Y0	9	SER
46	Y0	10	THR
46	Y0	11	ARG
46	Y0	36	ILE
46	Y0	41	ARG
46	Y0	64	ASP
46	Y0	66	VAL
46	Y0	74	ARG
47	Y1	21	ARG
47	Y1	30	VAL
47	Y1	46	LEU
47	Y1	50	ARG
47	Y1	51	VAL
47	Y1	56	GLN
47	Y1	62	VAL
47	Y1	67	ILE
47	Y1	78	LYS
47	Y1	80	LEU
47	Y1	82	LEU
47	Y1	83	GLU
47	Y1	91	LYS
48	Y2	7	ARG
48	Y2	9	GLN
48	Y2	16	LEU
48	Y2	24	LEU
48	Y2	53	LEU
48	Y2	62	THR
48	Y2	64	LEU
49	Y3	6	VAL
49	Y3	8	LEU
49	Y3	23	LEU
49	Y3	31	LEU
49	Y3	36	VAL
49	Y3	37	LEU
49	Y3	56	VAL
50	Y4	6	HIS
50	Y4	10	VAL
50	Y4	15	ILE
50	Y4	16	CYS
50	Y4	21	VAL
50	Y4	22	ILE
50	Y4	27	THR

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Mol	Chain	Res	Type
50	Y4	34	GLU
50	Y4	35	VAL
50	Y4	42	PHE
50	Y4	43	TYR
50	Y4	48	ARG
50	Y4	49	PHE
50	Y4	50	VAL
50	Y4	53	GLU
50	Y4	57	GLU
50	Y4	61	ARG
50	Y4	63	TYR
50	Y4	68	ARG
50	Y4	71	ARG
51	Y5	4	HIS
51	Y5	6	VAL
51	Y5	11	THR
51	Y5	29	THR
51	Y5	36	CYS
51	Y5	37	LYS
51	Y5	40	LYS
51	Y5	48	GLU
51	Y5	49	CYS
51	Y5	51	TYR
51	Y5	52	TYR
51	Y5	56	LYS
51	Y5	58	LEU
51	Y5	60	VAL
52	Y6	6	ARG
52	Y6	8	LYS
52	Y6	10	LEU
52	Y6	11	LEU
52	Y6	18	ARG
52	Y6	19	ARG
52	Y6	23	THR
52	Y6	27	LYS
52	Y6	30	THR
52	Y6	33	LYS
52	Y6	34	LEU
52	Y6	37	ARG
52	Y6	44	ARG
52	Y6	53	LYS
53	Y7	1	MET

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Mol	Chain	Res	Type
53	Y7	4	THR
53	Y7	8	ASN
53	Y7	9	ARG
53	Y7	14	LYS
53	Y7	46	VAL
53	Y7	47	ARG
54	Y8	16	ILE
54	Y8	27	THR
54	Y8	30	ARG
54	Y8	35	GLN
54	Y8	39	LYS
54	Y8	43	GLN
54	Y8	47	LYS
54	Y8	48	PHE
54	Y8	49	VAL
54	Y8	52	LYS
54	Y8	53	PRO
54	Y8	59	LYS
54	Y8	62	LEU
54	Y8	63	PRO
54	Y8	65	GLU
55	Y9	1	MET
55	Y9	17	ILE
55	Y9	26	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (92) such sidechains are listed below:

Mol	Chain	Res	Type
2	QB	94	ASN
2	QB	212	GLN
3	QC	139	GLN
3	QC	162	GLN
4	QD	45	GLN
7	QG	11	GLN
7	QG	64	GLN
7	QG	97	GLN
10	QJ	13	HIS
10	QJ	56	HIS
10	QJ	78	ASN
11	QK	93	GLN
11	QK	116	HIS
12	QL	9	GLN

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Mol	Chain	Res	Type
13	QM	106	ASN
15	QO	9	GLN
19	QS	47	HIS
19	QS	57	HIS
20	QT	90	GLN
28	RE	35	GLN
28	RE	55	ASN
28	RE	180	ASN
29	RF	69	HIS
29	RF	160	ASN
30	RG	40	ASN
30	RG	123	ASN
31	RH	143	GLN
31	RH	147	ASN
31	RH	158	HIS
32	RI	11	ASN
33	RN	130	HIS
36	RQ	123	HIS
38	RS	68	GLN
39	RT	2	ASN
39	RT	79	HIS
39	RT	84	GLN
40	RU	72	HIS
40	RU	81	HIS
44	RY	57	GLN
45	RZ	73	GLN
45	RZ	132	ASN
48	R2	43	GLN
48	R2	56	GLN
50	R4	6	HIS
51	R5	22	HIS
54	R8	33	ASN
55	R9	20	HIS
55	R9	29	ASN
55	R9	32	HIS
2	XB	40	HIS
2	XB	212	GLN
3	XC	102	ASN
3	XC	162	GLN
5	XE	65	ASN
5	XE	72	GLN
5	XE	141	GLN

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Mol	Chain	Res	Type
6	XF	27	GLN
7	XG	86	GLN
7	XG	106	GLN
7	XG	148	ASN
10	XJ	76	ASN
10	XJ	78	ASN
11	XK	38	ASN
12	XL	9	GLN
15	XO	50	HIS
15	XO	62	GLN
16	XP	13	HIS
27	YD	44	ASN
27	YD	87	ASN
27	YD	112	GLN
27	YD	164	GLN
27	YD	166	GLN
27	YD	198	ASN
28	YE	55	ASN
28	YE	180	ASN
30	YG	40	ASN
31	YH	143	GLN
31	YH	147	ASN
31	YH	158	HIS
33	YN	133	GLN
34	YO	90	GLN
35	YP	84	ASN
36	YQ	141	GLN
37	YR	71	GLN
42	YW	40	ASN
44	YY	57	GLN
45	YZ	151	HIS
48	Y2	9	GLN
48	Y2	38	GLN
48	Y2	46	GLN
54	Y8	33	ASN
55	Y9	29	ASN

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	QA	1497/1522 (98%)	277 (18%)	43 (2%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	XA	1498/1522 (98%)	294 (19%)	37 (2%)
22	QV	76/77 (98%)	19 (25%)	1 (1%)
22	XV	76/77 (98%)	14 (18%)	1 (1%)
23	QX	9/25 (36%)	1 (11%)	0
23	XX	9/25 (36%)	0	0
24	QY	14/17 (82%)	2 (14%)	0
24	XY	15/17 (88%)	1 (6%)	0
25	RA	2879/2915 (98%)	587 (20%)	58 (2%)
25	YA	2879/2915 (98%)	598 (20%)	52 (1%)
26	RB	119/122 (97%)	15 (12%)	2 (1%)
26	YB	119/122 (97%)	21 (17%)	1 (0%)
56	Z5	1/3 (33%)	0	0
56	Z6	1/3 (33%)	0	0
All	All	9192/9362 (98%)	1829 (19%)	195 (2%)

All (1829) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	QA	6	G
1	QA	9	G
1	QA	32	A
1	QA	39	G
1	QA	47	C
1	QA	48	C
1	QA	51	A
1	QA	64	G
1	QA	65	U
1	QA	66	G
1	QA	89	U
1	QA	91	C
1	QA	95	G
1	QA	101	A
1	QA	108	G
1	QA	116	A
1	QA	121	C
1	QA	129(A)	G
1	QA	144	G
1	QA	146	G
1	QA	147	G
1	QA	163	C
1	QA	169	C
1	QA	171	A

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Mol	Chain	Res	Type
1	QA	172	A
1	QA	173	U
1	QA	174	C
1	QA	182	U
1	QA	190	G
1	QA	191(A)	G
1	QA	195	A
1	QA	197	A
1	QA	209	U
1	QA	210	U
1	QA	216	G
1	QA	244	U
1	QA	245	C
1	QA	247	G
1	QA	251	G
1	QA	267	C
1	QA	270	A
1	QA	281	G
1	QA	289	G
1	QA	314	C
1	QA	321	A
1	QA	328	C
1	QA	329	A
1	QA	332	G
1	QA	344	A
1	QA	346	G
1	QA	347	G
1	QA	348	G
1	QA	352	C
1	QA	353	A
1	QA	354	G
1	QA	367	U
1	QA	372	C
1	QA	373	A
1	QA	389	A
1	QA	397	A
1	QA	398	C
1	QA	406	G
1	QA	411	A
1	QA	412	A
1	QA	413	G
1	QA	421	U

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Mol	Chain	Res	Type
1	QA	422	C
1	QA	423	G
1	QA	429	U
1	QA	430	A
1	QA	438	G
1	QA	442	C
1	QA	452	A
1	QA	466	C
1	QA	467	G
1	QA	482	A
1	QA	485	G
1	QA	486	U
1	QA	496	A
1	QA	497	U
1	QA	505	G
1	QA	509	A
1	QA	510	A
1	QA	511	C
1	QA	518	C
1	QA	521	G
1	QA	527	G
1	QA	531	U
1	QA	532	A
1	QA	533	A
1	QA	545	C
1	QA	547	A
1	QA	558	G
1	QA	559	A
1	QA	566	G
1	QA	568	G
1	QA	572	A
1	QA	573	A
1	QA	576	G
1	QA	577	G
1	QA	596	C
1	QA	614	A
1	QA	618	C
1	QA	630	G
1	QA	631	G
1	QA	632	A
1	QA	652	U
1	QA	653	A

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Mol	Chain	Res	Type
1	QA	661	G
1	QA	665	A
1	QA	686	U
1	QA	688	G
1	QA	701	C
1	QA	702	A
1	QA	703	G
1	QA	704	A
1	QA	721	G
1	QA	723	U
1	QA	724	G
1	QA	729	A
1	QA	731	G
1	QA	748	C
1	QA	749	C
1	QA	752	G
1	QA	753	A
1	QA	754	C
1	QA	755	G
1	QA	777	A
1	QA	786	G
1	QA	792	A
1	QA	793	U
1	QA	794	A
1	QA	813	U
1	QA	817	C
1	QA	819	A
1	QA	821	G
1	QA	828	A
1	QA	841	U
1	QA	842	C
1	QA	843	U
1	QA	848	C
1	QA	859	A
1	QA	870	U
1	QA	871	U
1	QA	872	A
1	QA	885	G
1	QA	902	G
1	QA	914	A
1	QA	927	G
1	QA	934	C

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Mol	Chain	Res	Type
1	QA	935	A
1	QA	960	U
1	QA	961	U
1	QA	968	A
1	QA	969	A
1	QA	971	G
1	QA	974	A
1	QA	975	A
1	QA	976	G
1	QA	977	A
1	QA	983	A
1	QA	991	U
1	QA	992	U
1	QA	993	G
1	QA	994	A
1	QA	1001	G
1	QA	1004	A
1	QA	1006	C
1	QA	1009	G
1	QA	1020	U
1	QA	1024	G
1	QA	1025	U
1	QA	1026	G
1	QA	1028	C
1	QA	1029	G
1	QA	1032(A)	G
1	QA	1036	G
1	QA	1040	U
1	QA	1046	A
1	QA	1053	G
1	QA	1064	G
1	QA	1065	U
1	QA	1066	C
1	QA	1081	G
1	QA	1094	G
1	QA	1095	U
1	QA	1101	A
1	QA	1124	G
1	QA	1125	U
1	QA	1126	U
1	QA	1130	A
1	QA	1131	G

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Mol	Chain	Res	Type
1	QA	1137	C
1	QA	1138	G
1	QA	1139	G
1	QA	1140	C
1	QA	1146	A
1	QA	1151	A
1	QA	1157	A
1	QA	1158	C
1	QA	1159	U
1	QA	1160	G
1	QA	1161	C
1	QA	1171	G
1	QA	1178	G
1	QA	1181	G
1	QA	1182	G
1	QA	1183	A
1	QA	1187	G
1	QA	1196	U
1	QA	1197	G
1	QA	1211	U
1	QA	1212	U
1	QA	1213	A
1	QA	1215	G
1	QA	1225	A
1	QA	1227	A
1	QA	1238	A
1	QA	1240	U
1	QA	1241	G
1	QA	1256	A
1	QA	1257	U
1	QA	1258	G
1	QA	1270	C
1	QA	1280	A
1	QA	1281	U
1	QA	1282	C
1	QA	1286	A
1	QA	1287	A
1	QA	1289	A
1	QA	1297	C
1	QA	1298	C
1	QA	1299	A
1	QA	1301	U

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Mol	Chain	Res	Type
1	QA	1302	U
1	QA	1303	C
1	QA	1305	G
1	QA	1312	G
1	QA	1319	A
1	QA	1320	C
1	QA	1321	C
1	QA	1322	C
1	QA	1323	G
1	QA	1331	G
1	QA	1335	C
1	QA	1336	C
1	QA	1337	G
1	QA	1338	G
1	QA	1346	A
1	QA	1347	G
1	QA	1348	U
1	QA	1353	G
1	QA	1362(A)	C
1	QA	1363	A
1	QA	1368	G
1	QA	1370	G
1	QA	1379	G
1	QA	1399	C
1	QA	1419	G
1	QA	1442	G
1	QA	1443	G
1	QA	1446	A
1	QA	1447	G
1	QA	1452	C
1	QA	1453	G
1	QA	1454	G
1	QA	1492	A
1	QA	1499	A
1	QA	1502	A
1	QA	1504	G
1	QA	1505	G
1	QA	1506	U
1	QA	1517	G
1	QA	1519	A
1	QA	1520	G
1	QA	1529	G

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Mol	Chain	Res	Type
1	QA	1530	G
22	QV	3	G
22	QV	4	G
22	QV	7	U
22	QV	12	C
22	QV	14	G
22	QV	15	C
22	QV	17	U
22	QV	18	G
22	QV	19	G
22	QV	21	A
22	QV	22	G
22	QV	47	U
22	QV	48	C
22	QV	53	G
22	QV	54	U
22	QV	59	A
22	QV	64	G
22	QV	67	C
22	QV	76	A
23	QX	19	A
24	QY	34	G
24	QY	36	U
25	RA	15	G
25	RA	34	C
25	RA	35	G
25	RA	46	C
25	RA	51	G
25	RA	55	G
25	RA	71	A
25	RA	72	U
25	RA	74	A
25	RA	75	G
25	RA	82	G
25	RA	91	A
25	RA	101	G
25	RA	102	G
25	RA	103	A
25	RA	118	A
25	RA	119	A
25	RA	120	U
25	RA	123	G

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Mol	Chain	Res	Type
25	RA	131	G
25	RA	138	G
25	RA	161	U
25	RA	177	G
25	RA	181	A
25	RA	196	A
25	RA	199	A
25	RA	215	G
25	RA	216	A
25	RA	221	A
25	RA	223	A
25	RA	228	A
25	RA	229	A
25	RA	230	U
25	RA	232	G
25	RA	238	C
25	RA	242	G
25	RA	243	U
25	RA	248	G
25	RA	249	C
25	RA	252	G
25	RA	265	A
25	RA	266	G
25	RA	269	U
25	RA	270(L)	U
25	RA	270(M)	U
25	RA	270(P)	C
25	RA	271(C)	U
25	RA	271	G
25	RA	275	G
25	RA	276	A
25	RA	277	C
25	RA	285	C
25	RA	286	C
25	RA	299	A
25	RA	311	A
25	RA	317	G
25	RA	323	G
25	RA	324	A
25	RA	327	G
25	RA	329	G
25	RA	330	A

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Mol	Chain	Res	Type
25	RA	333	G
25	RA	342	G
25	RA	346	A
25	RA	352	G
25	RA	363(F)	A
25	RA	364	C
25	RA	371	A
25	RA	372	G
25	RA	373	U
25	RA	386	G
25	RA	394	A
25	RA	395	U
25	RA	405	U
25	RA	411	G
25	RA	412	A
25	RA	428	A
25	RA	444	C
25	RA	448	U
25	RA	451	C
25	RA	455	C
25	RA	456	C
25	RA	457	A
25	RA	470	A
25	RA	481	G
25	RA	496	G
25	RA	504	U
25	RA	505	A
25	RA	508	G
25	RA	509	C
25	RA	512	G
25	RA	513	A
25	RA	527	C
25	RA	529	A
25	RA	531	C
25	RA	532	A
25	RA	533	G
25	RA	537	C
25	RA	539	G
25	RA	540	G
25	RA	546	C
25	RA	549	G
25	RA	554	U

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Mol	Chain	Res	Type
25	RA	556	G
25	RA	563	G
25	RA	573	G
25	RA	574	C
25	RA	575	A
25	RA	603	A
25	RA	604	G
25	RA	607	U
25	RA	609(A)	G
25	RA	614	U
25	RA	615	G
25	RA	617	G
25	RA	621	A
25	RA	627	A
25	RA	637	A
25	RA	638	G
25	RA	645	C
25	RA	646	A
25	RA	650	C
25	RA	651	G
25	RA	652	C
25	RA	653	A
25	RA	654	A
25	RA	654(A)	G
25	RA	668	G
25	RA	669	G
25	RA	686	G
25	RA	702	G
25	RA	717	G
25	RA	721	C
25	RA	722	A
25	RA	726	G
25	RA	730	C
25	RA	753	C
25	RA	764	A
25	RA	765	G
25	RA	775	G
25	RA	776	G
25	RA	782	A
25	RA	784	A
25	RA	785	G
25	RA	789	A

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Mol	Chain	Res	Type
25	RA	790	C
25	RA	792	G
25	RA	805	G
25	RA	812	C
25	RA	819	A
25	RA	827	U
25	RA	828	U
25	RA	831	G
25	RA	846	C
25	RA	847	U
25	RA	856	C
25	RA	857	C
25	RA	859	G
25	RA	860	U
25	RA	869	G
25	RA	882	G
25	RA	884	C
25	RA	885	C
25	RA	886	C
25	RA	887	A
25	RA	888	C
25	RA	889	C
25	RA	893	C
25	RA	896	A
25	RA	897	C
25	RA	899	A
25	RA	900	A
25	RA	901	A
25	RA	904	C
25	RA	906	G
25	RA	907	U
25	RA	910	A
25	RA	915	C
25	RA	917	A
25	RA	932	G
25	RA	933	A
25	RA	938	G
25	RA	941	A
25	RA	945	A
25	RA	946	G
25	RA	953	A
25	RA	961	C

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Mol	Chain	Res	Type
25	RA	974	G
25	RA	974(A)	C
25	RA	975	G
25	RA	980	A
25	RA	983	A
25	RA	996	A
25	RA	1003	G
25	RA	1011	G
25	RA	1012	U
25	RA	1013	C
25	RA	1015	G
25	RA	1017	G
25	RA	1020	A
25	RA	1023	U
25	RA	1025	G
25	RA	1026	U
25	RA	1027	A
25	RA	1033	U
25	RA	1044	G
25	RA	1045	A
25	RA	1046	A
25	RA	1050	A
25	RA	1055	G
25	RA	1057	A
25	RA	1059	G
25	RA	1060	U
25	RA	1061	U
25	RA	1065	U
25	RA	1066	U
25	RA	1067	A
25	RA	1068	G
25	RA	1071	G
25	RA	1073	A
25	RA	1077	A
25	RA	1078	U
25	RA	1079	C
25	RA	1080	C
25	RA	1082	U
25	RA	1083	U
25	RA	1084	A
25	RA	1085	A
25	RA	1086	A

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Mol	Chain	Res	Type
25	RA	1087	G
25	RA	1088	A
25	RA	1091	G
25	RA	1093	G
25	RA	1095	A
25	RA	1096	A
25	RA	1104	C
25	RA	1110	G
25	RA	1111	A
25	RA	1112	G
25	RA	1126	A
25	RA	1131	G
25	RA	1135	C
25	RA	1136	G
25	RA	1140	C
25	RA	1142	U
25	RA	1142(A)	A
25	RA	1151	G
25	RA	1173	G
25	RA	1174	A
25	RA	1175	U
25	RA	1176	G
25	RA	1179	C
25	RA	1180	C
25	RA	1183	G
25	RA	1195	G
25	RA	1204	A
25	RA	1205	U
25	RA	1206	G
25	RA	1210	A
25	RA	1211	U
25	RA	1220	A
25	RA	1238	G
25	RA	1240	U
25	RA	1248	G
25	RA	1253	A
25	RA	1256	G
25	RA	1265	A
25	RA	1271	G
25	RA	1272	A
25	RA	1276	A
25	RA	1288	U

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Mol	Chain	Res	Type
25	RA	1300	U
25	RA	1301	A
25	RA	1312	U
25	RA	1313	U
25	RA	1314	C
25	RA	1316	U
25	RA	1320	C
25	RA	1321	A
25	RA	1329	U
25	RA	1332	G
25	RA	1345	C
25	RA	1349	A
25	RA	1352	U
25	RA	1365	A
25	RA	1368	G
25	RA	1370	C
25	RA	1378	A
25	RA	1379	A
25	RA	1380	G
25	RA	1384	A
25	RA	1385	G
25	RA	1386	C
25	RA	1391	U
25	RA	1395	A
25	RA	1407	C
25	RA	1408	C
25	RA	1411	C
25	RA	1416	G
25	RA	1419	A
25	RA	1420	U
25	RA	1421	G
25	RA	1428	C
25	RA	1444(A)	A
25	RA	1445	C
25	RA	1449	A
25	RA	1449(A)	G
25	RA	1455	G
25	RA	1458	C
25	RA	1459	G
25	RA	1460	A
25	RA	1461	G
25	RA	1467	C

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Mol	Chain	Res	Type
25	RA	1471	A
25	RA	1480	G
25	RA	1482	U
25	RA	1483	G
25	RA	1485	G
25	RA	1486	A
25	RA	1493	C
25	RA	1494	A
25	RA	1496	A
25	RA	1497	U
25	RA	1504	C
25	RA	1505	C
25	RA	1506	C
25	RA	1507	A
25	RA	1508	A
25	RA	1510	A
25	RA	1513	C
25	RA	1514	U
25	RA	1515	C
25	RA	1519	G
25	RA	1522	G
25	RA	1533	C
25	RA	1534	G
25	RA	1535	U
25	RA	1536	A
25	RA	1537	C
25	RA	1538	G
25	RA	1543	A
25	RA	1545	A
25	RA	1554	A
25	RA	1558	A
25	RA	1559	G
25	RA	1566	A
25	RA	1569	A
25	RA	1578	U
25	RA	1579	A
25	RA	1580	A
25	RA	1581	G
25	RA	1585	C
25	RA	1586	A
25	RA	1598	C
25	RA	1607	C

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Mol	Chain	Res	Type
25	RA	1608	A
25	RA	1609	A
25	RA	1613	G
25	RA	1616	A
25	RA	1617	C
25	RA	1618	A
25	RA	1640	C
25	RA	1648	C
25	RA	1654	A
25	RA	1674	G
25	RA	1675	C
25	RA	1695	G
25	RA	1725	G
25	RA	1728	G
25	RA	1729	A
25	RA	1730	U
25	RA	1733	G
25	RA	1734	C
25	RA	1742	C
25	RA	1743	G
25	RA	1756	G
25	RA	1762	A
25	RA	1763	G
25	RA	1764	G
25	RA	1769	G
25	RA	1773	A
25	RA	1780	A
25	RA	1782	C
25	RA	1791	A
25	RA	1799	G
25	RA	1800	C
25	RA	1816	G
25	RA	1820	U
25	RA	1826	G
25	RA	1829	A
25	RA	1834	U
25	RA	1835	G
25	RA	1847	A
25	RA	1848	A
25	RA	1858	G
25	RA	1869	G
25	RA	1872	A

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Mol	Chain	Res	Type
25	RA	1878	G
25	RA	1882	C
25	RA	1885	A
25	RA	1888	G
25	RA	1889	A
25	RA	1906	G
25	RA	1913	A
25	RA	1914	C
25	RA	1929	G
25	RA	1931	U
25	RA	1934	C
25	RA	1936	A
25	RA	1938	A
25	RA	1939	U
25	RA	1955	U
25	RA	1963	U
25	RA	1967	C
25	RA	1969	A
25	RA	1970	A
25	RA	1971	A
25	RA	1972	A
25	RA	1982	C
25	RA	1991	U
25	RA	1992	G
25	RA	1993	U
25	RA	1996	C
25	RA	2021	C
25	RA	2023	G
25	RA	2031	A
25	RA	2032	G
25	RA	2033	A
25	RA	2043	C
25	RA	2055	C
25	RA	2056	G
25	RA	2059	A
25	RA	2060	A
25	RA	2061	G
25	RA	2062	A
25	RA	2069	G
25	RA	2099	U
25	RA	2108	C
25	RA	2111	C

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Mol	Chain	Res	Type
25	RA	2113	U
25	RA	2114	A
25	RA	2115	G
25	RA	2116	G
25	RA	2117	A
25	RA	2118	U
25	RA	2119	A
25	RA	2120	G
25	RA	2126	A
25	RA	2127	G
25	RA	2128	C
25	RA	2131	G
25	RA	2132	U
25	RA	2133	G
25	RA	2134	A
25	RA	2135	A
25	RA	2136	C
25	RA	2145	C
25	RA	2146	C
25	RA	2147	G
25	RA	2148	G
25	RA	2157	G
25	RA	2158	A
25	RA	2166	G
25	RA	2168	G
25	RA	2171	A
25	RA	2173	A
25	RA	2176	A
25	RA	2190	G
25	RA	2192	G
25	RA	2198	A
25	RA	2210	G
25	RA	2211	G
25	RA	2212	A
25	RA	2213	U
25	RA	2215	G
25	RA	2225	A
25	RA	2238	G
25	RA	2239	G
25	RA	2243	U
25	RA	2275	C
25	RA	2280	G

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Mol	Chain	Res	Type
25	RA	2283	C
25	RA	2287	A
25	RA	2288	A
25	RA	2298	A
25	RA	2300	G
25	RA	2307	G
25	RA	2308	G
25	RA	2311	A
25	RA	2312	U
25	RA	2319	G
25	RA	2320	A
25	RA	2325	G
25	RA	2335	A
25	RA	2336	A
25	RA	2345	G
25	RA	2346	A
25	RA	2347	C
25	RA	2350	C
25	RA	2354	G
25	RA	2379	G
25	RA	2383	G
25	RA	2385	C
25	RA	2392	A
25	RA	2402	C
25	RA	2403	C
25	RA	2406	U
25	RA	2410	G
25	RA	2411	A
25	RA	2423	U
25	RA	2424	C
25	RA	2425	A
25	RA	2429	G
25	RA	2430	A
25	RA	2435	A
25	RA	2439	A
25	RA	2440	C
25	RA	2441	C
25	RA	2448	A
25	RA	2469	A
25	RA	2470	G
25	RA	2475	C
25	RA	2482	G

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Mol	Chain	Res	Type
25	RA	2494	G
25	RA	2502	G
25	RA	2505	G
25	RA	2519	U
25	RA	2529	G
25	RA	2542	A
25	RA	2543	G
25	RA	2554	U
25	RA	2567	G
25	RA	2573	C
25	RA	2602	A
25	RA	2609	U
25	RA	2611	U
25	RA	2612	C
25	RA	2614	A
25	RA	2623	G
25	RA	2629	A
25	RA	2641	G
25	RA	2646	C
25	RA	2655	G
25	RA	2665	A
25	RA	2673	G
25	RA	2689	U
25	RA	2690	C
25	RA	2702	U
25	RA	2703	C
25	RA	2707	G
25	RA	2712	U
25	RA	2712(A)	A
25	RA	2713	A
25	RA	2714	G
25	RA	2726	U
25	RA	2733	A
25	RA	2734	A
25	RA	2748	A
25	RA	2752	C
25	RA	2758	A
25	RA	2759	G
25	RA	2761	G
25	RA	2764	A
25	RA	2765	A
25	RA	2766	G

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Mol	Chain	Res	Type
25	RA	2770	G
25	RA	2777	G
25	RA	2778	A
25	RA	2779	U
25	RA	2780	G
25	RA	2790	A
25	RA	2791	C
25	RA	2797	U
25	RA	2807	G
25	RA	2818	G
25	RA	2820	A
25	RA	2821	A
25	RA	2833	G
25	RA	2834	G
25	RA	2835	A
25	RA	2867	G
25	RA	2868	A
25	RA	2872	G
25	RA	2880	C
25	RA	2891	G
25	RA	2892	A
25	RA	2894	G
26	RB	9	G
26	RB	13	A
26	RB	15	A
26	RB	16	G
26	RB	22	U
26	RB	25	A
26	RB	27	C
26	RB	41	U
26	RB	42	C
26	RB	45	A
26	RB	56	G
26	RB	67	G
26	RB	73	A
26	RB	81	G
26	RB	109	G
1	XA	6	G
1	XA	7	G
1	XA	9	G
1	XA	32	A
1	XA	39	G

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Mol	Chain	Res	Type
1	XA	47	C
1	XA	48	C
1	XA	50	A
1	XA	51	A
1	XA	61	G
1	XA	64	G
1	XA	65	U
1	XA	66	G
1	XA	76	G
1	XA	78	G
1	XA	79	G
1	XA	81	G
1	XA	88	C
1	XA	89	U
1	XA	90	C
1	XA	91	C
1	XA	92	G
1	XA	95	G
1	XA	101	A
1	XA	108	G
1	XA	116	A
1	XA	121	C
1	XA	129(A)	G
1	XA	130	A
1	XA	144	G
1	XA	147	G
1	XA	163	C
1	XA	169	C
1	XA	172	A
1	XA	173	U
1	XA	174	C
1	XA	190	G
1	XA	191(A)	G
1	XA	195	A
1	XA	197	A
1	XA	201	C
1	XA	209	U
1	XA	216	G
1	XA	222	U
1	XA	243	A
1	XA	244	U
1	XA	247	G

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Mol	Chain	Res	Type
1	XA	251	G
1	XA	266	G
1	XA	267	C
1	XA	270	A
1	XA	281	G
1	XA	289	G
1	XA	321	A
1	XA	328	C
1	XA	329	A
1	XA	332	G
1	XA	345	C
1	XA	346	G
1	XA	347	G
1	XA	348	G
1	XA	352	C
1	XA	353	A
1	XA	354	G
1	XA	356	A
1	XA	367	U
1	XA	372	C
1	XA	373	A
1	XA	381	C
1	XA	384	G
1	XA	388	G
1	XA	389	A
1	XA	397	A
1	XA	398	C
1	XA	406	G
1	XA	412	A
1	XA	413	G
1	XA	414	A
1	XA	421	U
1	XA	422	C
1	XA	423	G
1	XA	424	G
1	XA	429	U
1	XA	430	A
1	XA	435	C
1	XA	439	A
1	XA	442	C
1	XA	466	C
1	XA	467	G

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Mol	Chain	Res	Type
1	XA	482	A
1	XA	485	G
1	XA	486	U
1	XA	487	A
1	XA	496	A
1	XA	497	U
1	XA	509	A
1	XA	510	A
1	XA	511	C
1	XA	518	C
1	XA	521	G
1	XA	527	G
1	XA	531	U
1	XA	532	A
1	XA	533	A
1	XA	547	A
1	XA	548	G
1	XA	559	A
1	XA	561	U
1	XA	564	C
1	XA	568	G
1	XA	572	A
1	XA	573	A
1	XA	576	G
1	XA	577	G
1	XA	579	G
1	XA	596	C
1	XA	619	U
1	XA	630	G
1	XA	631	G
1	XA	632	A
1	XA	633	G
1	XA	653	A
1	XA	661	G
1	XA	665	A
1	XA	671	G
1	XA	687	A
1	XA	688	G
1	XA	704	A
1	XA	721	G
1	XA	723	U
1	XA	724	G

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Mol	Chain	Res	Type
1	XA	731	G
1	XA	734	G
1	XA	748	C
1	XA	755	G
1	XA	774	G
1	XA	777	A
1	XA	792	A
1	XA	793	U
1	XA	794	A
1	XA	813	U
1	XA	816	A
1	XA	817	C
1	XA	818	G
1	XA	828	A
1	XA	841	U
1	XA	842	C
1	XA	843	U
1	XA	848	C
1	XA	859	A
1	XA	864	A
1	XA	871	U
1	XA	872	A
1	XA	881	G
1	XA	902	G
1	XA	914	A
1	XA	927	G
1	XA	934	C
1	XA	935	A
1	XA	936	C
1	XA	948	C
1	XA	960	U
1	XA	966	G
1	XA	968	A
1	XA	969	A
1	XA	971	G
1	XA	974	A
1	XA	975	A
1	XA	976	G
1	XA	977	A
1	XA	983	A
1	XA	991	U
1	XA	992	U

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Mol	Chain	Res	Type
1	XA	993	G
1	XA	1004	A
1	XA	1006	C
1	XA	1008	C
1	XA	1021	G
1	XA	1024	G
1	XA	1025	U
1	XA	1028	C
1	XA	1029	G
1	XA	1032(A)	G
1	XA	1033	G
1	XA	1036	G
1	XA	1040	U
1	XA	1042	G
1	XA	1050	G
1	XA	1053	G
1	XA	1054	C
1	XA	1055	A
1	XA	1081	G
1	XA	1085	U
1	XA	1094	G
1	XA	1095	U
1	XA	1096	C
1	XA	1101	A
1	XA	1124	G
1	XA	1125	U
1	XA	1126	U
1	XA	1127	G
1	XA	1130	A
1	XA	1131	G
1	XA	1136	U
1	XA	1137	C
1	XA	1138	G
1	XA	1139	G
1	XA	1145	C
1	XA	1146	A
1	XA	1157	A
1	XA	1158	C
1	XA	1159	U
1	XA	1160	G
1	XA	1162	C
1	XA	1171	G

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Mol	Chain	Res	Type
1	XA	1177	G
1	XA	1178	G
1	XA	1181	G
1	XA	1182	G
1	XA	1183	A
1	XA	1184	G
1	XA	1190	G
1	XA	1195	C
1	XA	1196	U
1	XA	1197	G
1	XA	1200	C
1	XA	1201	A
1	XA	1202	G
1	XA	1211	U
1	XA	1212	U
1	XA	1213	A
1	XA	1225	A
1	XA	1226	C
1	XA	1236	A
1	XA	1238	A
1	XA	1240	U
1	XA	1241	G
1	XA	1256	A
1	XA	1257	U
1	XA	1258	G
1	XA	1260	C
1	XA	1263	C
1	XA	1270	C
1	XA	1272	G
1	XA	1273	G
1	XA	1280	A
1	XA	1281	U
1	XA	1282	C
1	XA	1286	A
1	XA	1287	A
1	XA	1297	C
1	XA	1298	C
1	XA	1300	G
1	XA	1301	U
1	XA	1302	U
1	XA	1305	G
1	XA	1317	C

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Mol	Chain	Res	Type
1	XA	1319	A
1	XA	1320	C
1	XA	1321	C
1	XA	1322	C
1	XA	1323	G
1	XA	1331	G
1	XA	1336	C
1	XA	1337	G
1	XA	1346	A
1	XA	1347	G
1	XA	1348	U
1	XA	1353	G
1	XA	1362(A)	C
1	XA	1363	A
1	XA	1364	U
1	XA	1370	G
1	XA	1419	G
1	XA	1437	C
1	XA	1438	G
1	XA	1442	G
1	XA	1446	A
1	XA	1447	G
1	XA	1452	C
1	XA	1453	G
1	XA	1487	G
1	XA	1492	A
1	XA	1497	G
1	XA	1499	A
1	XA	1502	A
1	XA	1503	A
1	XA	1506	U
1	XA	1517	G
1	XA	1519	A
1	XA	1520	G
1	XA	1529	G
1	XA	1530	G
1	XA	1531	A
22	XV	3	G
22	XV	4	G
22	XV	17	U
22	XV	18	G
22	XV	19	G

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Mol	Chain	Res	Type
22	XV	21	A
22	XV	25	C
22	XV	47	U
22	XV	48	C
22	XV	52	G
22	XV	54	U
22	XV	59	A
22	XV	67	C
22	XV	76	A
24	XY	36	U
25	YA	9	U
25	YA	15	G
25	YA	27	G
25	YA	34	C
25	YA	35	G
25	YA	46	C
25	YA	55	G
25	YA	71	A
25	YA	72	U
25	YA	74	A
25	YA	75	G
25	YA	96	G
25	YA	98	G
25	YA	99	U
25	YA	101	G
25	YA	102	G
25	YA	103	A
25	YA	118	A
25	YA	119	A
25	YA	120	U
25	YA	125	G
25	YA	131	G
25	YA	161	U
25	YA	181	A
25	YA	188	G
25	YA	196	A
25	YA	199	A
25	YA	214	G
25	YA	215	G
25	YA	216	A
25	YA	221	A
25	YA	222	A

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Mol	Chain	Res	Type
25	YA	223	A
25	YA	226	G
25	YA	228	A
25	YA	229	A
25	YA	230	U
25	YA	232	G
25	YA	242	G
25	YA	243	U
25	YA	248	G
25	YA	252	G
25	YA	265	A
25	YA	266	G
25	YA	269	U
25	YA	270(L)	U
25	YA	270(M)	U
25	YA	270(N)	G
25	YA	270(P)	C
25	YA	271(A)	C
25	YA	271(B)	G
25	YA	271(C)	U
25	YA	271	G
25	YA	274	G
25	YA	275	G
25	YA	276	A
25	YA	278	A
25	YA	279	C
25	YA	299	A
25	YA	311	A
25	YA	316	C
25	YA	323	G
25	YA	324	A
25	YA	329	G
25	YA	330	A
25	YA	332	A
25	YA	342	G
25	YA	345	A
25	YA	352	G
25	YA	363	G
25	YA	364	C
25	YA	371	A
25	YA	372	G
25	YA	386	G

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Mol	Chain	Res	Type
25	YA	387	U
25	YA	396	G
25	YA	405	U
25	YA	406	G
25	YA	411	G
25	YA	428	A
25	YA	443	A
25	YA	444	C
25	YA	448	U
25	YA	457	A
25	YA	470	A
25	YA	481	G
25	YA	496	G
25	YA	503	A
25	YA	504	U
25	YA	505	A
25	YA	508	G
25	YA	509	C
25	YA	512	G
25	YA	518	G
25	YA	531	C
25	YA	532	A
25	YA	533	G
25	YA	537	C
25	YA	539	G
25	YA	540	G
25	YA	546	C
25	YA	547	A
25	YA	549	G
25	YA	563	G
25	YA	571	A
25	YA	573	G
25	YA	575	A
25	YA	586	A
25	YA	588	U
25	YA	603	A
25	YA	604	G
25	YA	607	U
25	YA	614	U
25	YA	615	G
25	YA	616	A
25	YA	617	G

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Mol	Chain	Res	Type
25	YA	622	G
25	YA	627	A
25	YA	637	A
25	YA	638	G
25	YA	645	C
25	YA	646	A
25	YA	651	G
25	YA	654(A)	G
25	YA	654(B)	C
25	YA	654(V)	A
25	YA	668	G
25	YA	670	A
25	YA	686	G
25	YA	702	G
25	YA	717	G
25	YA	722	A
25	YA	726	G
25	YA	730	C
25	YA	734	A
25	YA	747	U
25	YA	753	C
25	YA	764	A
25	YA	775	G
25	YA	776	G
25	YA	782	A
25	YA	784	A
25	YA	785	G
25	YA	789	A
25	YA	790	C
25	YA	792	G
25	YA	805	G
25	YA	812	C
25	YA	819	A
25	YA	827	U
25	YA	828	U
25	YA	831	G
25	YA	847	U
25	YA	856	C
25	YA	857	C
25	YA	860	U
25	YA	865	C
25	YA	866	A

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Mol	Chain	Res	Type
25	YA	881	G
25	YA	882	G
25	YA	884	C
25	YA	885	C
25	YA	886	C
25	YA	888	C
25	YA	889	C
25	YA	892	G
25	YA	896	A
25	YA	898	C
25	YA	899	A
25	YA	900	A
25	YA	901	A
25	YA	906	G
25	YA	907	U
25	YA	910	A
25	YA	914	C
25	YA	915	C
25	YA	917	A
25	YA	918	A
25	YA	926	A
25	YA	932	G
25	YA	938	G
25	YA	941	A
25	YA	944	G
25	YA	945	A
25	YA	946	G
25	YA	961	C
25	YA	974	G
25	YA	974(A)	C
25	YA	975	G
25	YA	980	A
25	YA	983	A
25	YA	990	A
25	YA	991	C
25	YA	996	A
25	YA	1003	G
25	YA	1005	C
25	YA	1011	G
25	YA	1012	U
25	YA	1013	C
25	YA	1016	G

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Mol	Chain	Res	Type
25	YA	1017	G
25	YA	1022	G
25	YA	1023	U
25	YA	1025	G
25	YA	1026	U
25	YA	1027	A
25	YA	1033	U
25	YA	1045	A
25	YA	1046	A
25	YA	1050	A
25	YA	1054	A
25	YA	1055	G
25	YA	1059	G
25	YA	1060	U
25	YA	1061	U
25	YA	1066	U
25	YA	1067	A
25	YA	1068	G
25	YA	1071	G
25	YA	1076	C
25	YA	1077	A
25	YA	1078	U
25	YA	1079	C
25	YA	1082	U
25	YA	1083	U
25	YA	1084	A
25	YA	1085	A
25	YA	1086	A
25	YA	1088	A
25	YA	1089	G
25	YA	1090	U
25	YA	1095	A
25	YA	1096	A
25	YA	1097	U
25	YA	1103	A
25	YA	1104	C
25	YA	1105	U
25	YA	1110	G
25	YA	1111	A
25	YA	1122	G
25	YA	1131	G
25	YA	1135	C

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Mol	Chain	Res	Type
25	YA	1136	G
25	YA	1139	G
25	YA	1142	U
25	YA	1142(A)	A
25	YA	1151	G
25	YA	1170	G
25	YA	1173	G
25	YA	1174	A
25	YA	1175	U
25	YA	1176	G
25	YA	1178	C
25	YA	1179	C
25	YA	1180	C
25	YA	1195	G
25	YA	1204	A
25	YA	1205	U
25	YA	1210	A
25	YA	1211	U
25	YA	1220	A
25	YA	1221	C
25	YA	1236	G
25	YA	1238	G
25	YA	1250	G
25	YA	1252	G
25	YA	1253	A
25	YA	1256	G
25	YA	1265	A
25	YA	1271	G
25	YA	1272	A
25	YA	1273	U
25	YA	1275	A
25	YA	1287	A
25	YA	1300	U
25	YA	1301	A
25	YA	1302	A
25	YA	1313	U
25	YA	1319	G
25	YA	1321	A
25	YA	1329	U
25	YA	1345	C
25	YA	1349	A
25	YA	1352	U

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Mol	Chain	Res	Type
25	YA	1358	G
25	YA	1359	A
25	YA	1360	A
25	YA	1365	A
25	YA	1368	G
25	YA	1379	A
25	YA	1384	A
25	YA	1385	G
25	YA	1386	C
25	YA	1389	G
25	YA	1390	U
25	YA	1392	A
25	YA	1407	C
25	YA	1408	C
25	YA	1411	C
25	YA	1416	G
25	YA	1417	C
25	YA	1419	A
25	YA	1420	U
25	YA	1421	G
25	YA	1427	A
25	YA	1428	C
25	YA	1429	G
25	YA	1444(A)	A
25	YA	1445	C
25	YA	1449	A
25	YA	1449(A)	G
25	YA	1455	G
25	YA	1458	C
25	YA	1460	A
25	YA	1461	G
25	YA	1464	C
25	YA	1467	C
25	YA	1471	A
25	YA	1482	U
25	YA	1483	G
25	YA	1485	G
25	YA	1487	G
25	YA	1493	C
25	YA	1497	U
25	YA	1507	A
25	YA	1508	A

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Mol	Chain	Res	Type
25	YA	1510	A
25	YA	1511	A
25	YA	1514	U
25	YA	1515	C
25	YA	1522	G
25	YA	1534	G
25	YA	1535	U
25	YA	1536	A
25	YA	1537	C
25	YA	1543	A
25	YA	1544	C
25	YA	1545	A
25	YA	1554	A
25	YA	1558	A
25	YA	1559	G
25	YA	1566	A
25	YA	1569	A
25	YA	1578	U
25	YA	1579	A
25	YA	1581	G
25	YA	1585	C
25	YA	1586	A
25	YA	1587	A
25	YA	1591	G
25	YA	1598	C
25	YA	1608	A
25	YA	1609	A
25	YA	1617	C
25	YA	1618	A
25	YA	1640	C
25	YA	1647	G
25	YA	1648	C
25	YA	1654	A
25	YA	1668	A
25	YA	1674	G
25	YA	1694	C
25	YA	1695	G
25	YA	1700	A
25	YA	1701	A
25	YA	1725	G
25	YA	1728	G
25	YA	1729	A

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Mol	Chain	Res	Type
25	YA	1730	U
25	YA	1731	G
25	YA	1732	A
25	YA	1733	G
25	YA	1742	C
25	YA	1743	G
25	YA	1750	G
25	YA	1754	C
25	YA	1756	G
25	YA	1759	A
25	YA	1762	A
25	YA	1763	G
25	YA	1764	G
25	YA	1773	A
25	YA	1780	A
25	YA	1787	A
25	YA	1791	A
25	YA	1799	G
25	YA	1800	C
25	YA	1801	G
25	YA	1811	G
25	YA	1816	G
25	YA	1820	U
25	YA	1829	A
25	YA	1835	G
25	YA	1847	A
25	YA	1858	G
25	YA	1869	G
25	YA	1870	C
25	YA	1872	A
25	YA	1878	G
25	YA	1882	C
25	YA	1889	A
25	YA	1896	G
25	YA	1899	G
25	YA	1906	G
25	YA	1913	A
25	YA	1919	A
25	YA	1929	G
25	YA	1930	G
25	YA	1931	U
25	YA	1938	A

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Mol	Chain	Res	Type
25	YA	1939	U
25	YA	1955	U
25	YA	1956	U
25	YA	1963	U
25	YA	1965	C
25	YA	1967	C
25	YA	1969	A
25	YA	1970	A
25	YA	1971	A
25	YA	1972	A
25	YA	1982	C
25	YA	1991	U
25	YA	1993	U
25	YA	2020	A
25	YA	2021	C
25	YA	2023	G
25	YA	2031	A
25	YA	2033	A
25	YA	2034	U
25	YA	2043	C
25	YA	2055	C
25	YA	2056	G
25	YA	2059	A
25	YA	2060	A
25	YA	2061	G
25	YA	2062	A
25	YA	2063	C
25	YA	2069	G
25	YA	2089	U
25	YA	2093	G
25	YA	2099	U
25	YA	2100	G
25	YA	2111	C
25	YA	2112	G
25	YA	2113	U
25	YA	2114	A
25	YA	2115	G
25	YA	2116	G
25	YA	2117	A
25	YA	2120	G
25	YA	2122	U
25	YA	2126	A

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Mol	Chain	Res	Type
25	YA	2127	G
25	YA	2128	C
25	YA	2131	G
25	YA	2132	U
25	YA	2133	G
25	YA	2136	C
25	YA	2146	C
25	YA	2148	G
25	YA	2157	G
25	YA	2158	A
25	YA	2166	G
25	YA	2167	U
25	YA	2168	G
25	YA	2173	A
25	YA	2176	A
25	YA	2181	G
25	YA	2189	U
25	YA	2190	G
25	YA	2192	G
25	YA	2193	G
25	YA	2198	A
25	YA	2210	G
25	YA	2211	G
25	YA	2212	A
25	YA	2215	G
25	YA	2225	A
25	YA	2238	G
25	YA	2239	G
25	YA	2243	U
25	YA	2244	U
25	YA	2275	C
25	YA	2278	A
25	YA	2283	C
25	YA	2287	A
25	YA	2288	A
25	YA	2307	G
25	YA	2308	G
25	YA	2311	A
25	YA	2312	U
25	YA	2319	G
25	YA	2320	A
25	YA	2325	G

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Mol	Chain	Res	Type
25	YA	2334	G
25	YA	2346	A
25	YA	2347	C
25	YA	2350	C
25	YA	2358	G
25	YA	2377	A
25	YA	2383	G
25	YA	2385	C
25	YA	2392	A
25	YA	2394	C
25	YA	2402	C
25	YA	2403	C
25	YA	2406	U
25	YA	2410	G
25	YA	2423	U
25	YA	2424	C
25	YA	2425	A
25	YA	2427	C
25	YA	2429	G
25	YA	2430	A
25	YA	2435	A
25	YA	2439	A
25	YA	2440	C
25	YA	2441	C
25	YA	2448	A
25	YA	2450	A
25	YA	2469	A
25	YA	2470	G
25	YA	2471	C
25	YA	2475	C
25	YA	2484	G
25	YA	2494	G
25	YA	2498	C
25	YA	2502	G
25	YA	2505	G
25	YA	2518	A
25	YA	2529	G
25	YA	2542	A
25	YA	2554	U
25	YA	2558	C
25	YA	2567	G
25	YA	2601	C

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Mol	Chain	Res	Type
25	YA	2602	A
25	YA	2609	U
25	YA	2610	C
25	YA	2611	U
25	YA	2612	C
25	YA	2629	A
25	YA	2632	A
25	YA	2655	G
25	YA	2665	A
25	YA	2666	C
25	YA	2673	G
25	YA	2682	U
25	YA	2689	U
25	YA	2690	C
25	YA	2691	C
25	YA	2702	U
25	YA	2703	C
25	YA	2707	G
25	YA	2712	U
25	YA	2712(A)	A
25	YA	2713	A
25	YA	2714	G
25	YA	2718	G
25	YA	2726	U
25	YA	2733	A
25	YA	2742	C
25	YA	2744	G
25	YA	2752	C
25	YA	2758	A
25	YA	2761	G
25	YA	2764	A
25	YA	2765	A
25	YA	2766	G
25	YA	2770	G
25	YA	2777	G
25	YA	2778	A
25	YA	2779	U
25	YA	2780	G
25	YA	2789	C
25	YA	2790	A
25	YA	2791	C
25	YA	2794	C

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Mol	Chain	Res	Type
25	YA	2797	U
25	YA	2807	G
25	YA	2808	U
25	YA	2818	G
25	YA	2820	A
25	YA	2821	A
25	YA	2830	G
25	YA	2833	G
25	YA	2834	G
25	YA	2835	A
25	YA	2836	U
25	YA	2847	U
25	YA	2867	G
25	YA	2868	A
25	YA	2872	G
25	YA	2873	A
25	YA	2880	C
25	YA	2892	A
25	YA	2893	G
25	YA	2894	G
26	YB	8	U
26	YB	9	G
26	YB	13	A
26	YB	15	A
26	YB	16	G
26	YB	22	U
26	YB	25	A
26	YB	32	C
26	YB	40	U
26	YB	41	U
26	YB	42	C
26	YB	45	A
26	YB	52	A
26	YB	53	A
26	YB	56	G
26	YB	67	G
26	YB	73	A
26	YB	108	C
26	YB	109	G
26	YB	115	G
26	YB	116	G

All (195) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	QA	5	U
1	QA	31	G
1	QA	64	G
1	QA	115	G
1	QA	181	G
1	QA	189	U
1	QA	243	A
1	QA	244	U
1	QA	250	A
1	QA	266	G
1	QA	328	C
1	QA	410	G
1	QA	412	A
1	QA	428	G
1	QA	429	U
1	QA	481	G
1	QA	484	G
1	QA	485	G
1	QA	509	A
1	QA	687	A
1	QA	701	C
1	QA	703	G
1	QA	753	A
1	QA	792	A
1	QA	812	C
1	QA	913	A
1	QA	960	U
1	QA	991	U
1	QA	992	U
1	QA	1025	U
1	QA	1027	C
1	QA	1064	G
1	QA	1065	U
1	QA	1280	A
1	QA	1285	A
1	QA	1297	C
1	QA	1301	U
1	QA	1336	C
1	QA	1346	A
1	QA	1347	G
1	QA	1446	A
1	QA	1498	U
1	QA	1528	U

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Mol	Chain	Res	Type
22	QV	53	G
25	RA	74	A
25	RA	99	U
25	RA	102	G
25	RA	222	A
25	RA	227	A
25	RA	229	A
25	RA	242	G
25	RA	271(B)	G
25	RA	271(C)	U
25	RA	345	A
25	RA	370	G
25	RA	372	G
25	RA	404	C
25	RA	503	A
25	RA	508	G
25	RA	512	G
25	RA	637	A
25	RA	752	A
25	RA	764	A
25	RA	774	A
25	RA	846	C
25	RA	856	C
25	RA	859	G
25	RA	974(A)	C
25	RA	1022	G
25	RA	1026	U
25	RA	1045	A
25	RA	1078	U
25	RA	1085	A
25	RA	1130	U
25	RA	1178	C
25	RA	1204	A
25	RA	1210	A
25	RA	1275	A
25	RA	1312	U
25	RA	1427	A
25	RA	1460	A
25	RA	1558	A
25	RA	1653	G
25	RA	1694	C
25	RA	1799	G

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Mol	Chain	Res	Type
25	RA	1819	A
25	RA	1930	G
25	RA	1992	G
25	RA	2060	A
25	RA	2126	A
25	RA	2405	G
25	RA	2439	A
25	RA	2481	G
25	RA	2518	A
25	RA	2566	A
25	RA	2610	C
25	RA	2689	U
25	RA	2712	U
25	RA	2726	U
25	RA	2776	A
25	RA	2832	U
25	RA	2867	G
26	RB	24	G
26	RB	66	A
1	XA	31	G
1	XA	60	A
1	XA	64	G
1	XA	78	G
1	XA	89	U
1	XA	115	G
1	XA	190	G
1	XA	243	A
1	XA	250	A
1	XA	266	G
1	XA	328	C
1	XA	345	C
1	XA	353	A
1	XA	388	G
1	XA	412	A
1	XA	429	U
1	XA	481	G
1	XA	484	G
1	XA	485	G
1	XA	509	A
1	XA	532	A
1	XA	560	U
1	XA	687	A

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Mol	Chain	Res	Type
1	XA	703	G
1	XA	812	C
1	XA	913	A
1	XA	991	U
1	XA	992	U
1	XA	1027	C
1	XA	1054	C
1	XA	1200	C
1	XA	1285	A
1	XA	1301	U
1	XA	1336	C
1	XA	1347	G
1	XA	1446	A
1	XA	1498	U
22	XV	53	G
25	YA	99	U
25	YA	102	G
25	YA	195	A
25	YA	196	A
25	YA	221	A
25	YA	222	A
25	YA	229	A
25	YA	242	G
25	YA	271(B)	G
25	YA	278	A
25	YA	404	C
25	YA	503	A
25	YA	508	G
25	YA	587	C
25	YA	637	A
25	YA	654	A
25	YA	752	A
25	YA	846	C
25	YA	856	C
25	YA	859	G
25	YA	974(A)	C
25	YA	1012	U
25	YA	1022	G
25	YA	1026	U
25	YA	1045	A
25	YA	1078	U
25	YA	1085	A

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Mol	Chain	Res	Type
25	YA	1109	C
25	YA	1130	U
25	YA	1178	C
25	YA	1204	A
25	YA	1210	A
25	YA	1359	A
25	YA	1427	A
25	YA	1558	A
25	YA	1653	G
25	YA	1694	C
25	YA	1799	G
25	YA	1819	A
25	YA	1930	G
25	YA	1955	U
25	YA	1992	G
25	YA	2126	A
25	YA	2439	A
25	YA	2566	A
25	YA	2610	C
25	YA	2681	C
25	YA	2689	U
25	YA	2712	U
25	YA	2776	A
25	YA	2832	U
25	YA	2867	G
26	YB	66	A

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
56	PPU	Z6	101	25,56	37,40,41	2.58	8 (21%)	47,57,60	2.56	15 (31%)
56	PPU	Z5	101	25,56	37,40,41	2.58	8 (21%)	47,57,60	2.56	15 (31%)

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
56	PPU	Z6	101	25,56	-	2/25/43/44	0/4/4/4
56	PPU	Z5	101	25,56	-	2/25/43/44	0/4/4/4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
56	Z5	101	PPU	O-C	9.65	1.41	1.23
56	Z6	101	PPU	O-C	9.61	1.41	1.23
56	Z5	101	PPU	C9-N6	-6.20	1.32	1.45
56	Z6	101	PPU	C9-N6	-6.18	1.32	1.45
56	Z6	101	PPU	C10-N6	-5.89	1.32	1.45
56	Z5	101	PPU	C10-N6	-5.87	1.32	1.45
56	Z5	101	PPU	C-N3'	5.86	1.46	1.34
56	Z6	101	PPU	C-N3'	5.83	1.46	1.34
56	Z6	101	PPU	C5-N7	-3.60	1.32	1.39
56	Z5	101	PPU	C5-N7	-3.57	1.32	1.39
56	Z5	101	PPU	C8-N9	-3.19	1.32	1.37
56	Z6	101	PPU	C8-N9	-3.16	1.32	1.37
56	Z6	101	PPU	C1'-N9	2.74	1.53	1.46
56	Z5	101	PPU	C1'-N9	2.74	1.53	1.46
56	Z6	101	PPU	C4-N9	-2.24	1.33	1.37
56	Z5	101	PPU	C4-N9	-2.23	1.33	1.37

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
56	Z5	101	PPU	C3'-N3'-C	-8.53	110.22	123.20
56	Z6	101	PPU	C3'-N3'-C	-8.52	110.24	123.20
56	Z6	101	PPU	C2'-C1'-N9	-5.93	98.57	113.30
56	Z5	101	PPU	C2'-C1'-N9	-5.93	98.57	113.30
56	Z6	101	PPU	C5-C4-N3	-5.02	119.80	126.72
56	Z5	101	PPU	C5-C4-N3	-5.01	119.82	126.72
56	Z5	101	PPU	N3-C2-N1	-4.71	121.45	128.58
56	Z6	101	PPU	N3-C2-N1	-4.69	121.48	128.58
56	Z6	101	PPU	C4-C5-N7	-4.11	105.89	110.58
56	Z5	101	PPU	C4-C5-N7	-4.07	105.92	110.58
56	Z5	101	PPU	CA-C-N3'	4.07	121.72	116.21
56	Z6	101	PPU	CA-C-N3'	4.05	121.70	116.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
56	Z6	101	PPU	C2-N1-C6	4.02	121.64	111.83
56	Z5	101	PPU	C2-N1-C6	4.02	121.64	111.83
56	Z5	101	PPU	C2-N3-C4	3.68	120.81	111.83
56	Z6	101	PPU	C2-N3-C4	3.68	120.81	111.83
56	Z6	101	PPU	C5-N7-C8	3.63	109.16	103.45
56	Z5	101	PPU	C5-N7-C8	3.59	109.09	103.45
56	Z6	101	PPU	N9-C8-N7	-3.46	109.03	113.94
56	Z5	101	PPU	N9-C8-N7	-3.42	109.08	113.94
56	Z5	101	PPU	CM-OC-CZ	-3.42	110.17	117.50
56	Z6	101	PPU	CM-OC-CZ	-3.42	110.17	117.50
56	Z5	101	PPU	O4'-C1'-N9	-3.01	102.31	108.09
56	Z6	101	PPU	O4'-C1'-N9	-2.98	102.36	108.09
56	Z6	101	PPU	N3-C4-N9	2.87	132.04	127.17
56	Z5	101	PPU	N3-C4-N9	2.85	132.01	127.17
56	Z5	101	PPU	C4-C5-C6	2.49	118.49	115.91
56	Z6	101	PPU	C4-C5-C6	2.47	118.46	115.91
56	Z5	101	PPU	C5-C4-N9	2.16	108.16	105.81
56	Z6	101	PPU	C5-C4-N9	2.15	108.16	105.81

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
56	Z5	101	PPU	O-C-CA-N
56	Z6	101	PPU	O-C-CA-N
56	Z5	101	PPU	N3'-C-CA-N
56	Z6	101	PPU	N3'-C-CA-N

There are no ring outliers.

2 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
56	Z6	101	PPU	11	0
56	Z5	101	PPU	8	0

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 902 ligands modelled in this entry, 900 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	$\#\|Z\| > 2$
58	PAR	QA	1696	-	44,45,45	1.41	8 (18%)	63,67,67	1.27	6 (9%)
58	PAR	XA	1715	-	44,45,45	1.42	6 (13%)	63,67,67	1.21	4 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
58	PAR	QA	1696	-	-	4/18/94/94	0/4/4/4
58	PAR	XA	1715	-	-	6/18/94/94	0/4/4/4

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
58	QA	1696	PAR	C52-C42	3.32	1.59	1.52
58	XA	1715	PAR	C52-C42	3.14	1.58	1.52
58	XA	1715	PAR	O54-C14	3.06	1.49	1.41
58	QA	1696	PAR	O54-C14	2.97	1.49	1.41
58	QA	1696	PAR	C11-C21	2.73	1.57	1.52
58	XA	1715	PAR	O51-C11	2.65	1.48	1.41
58	XA	1715	PAR	C14-C24	2.64	1.57	1.52
58	QA	1696	PAR	O51-C11	2.58	1.48	1.41
58	XA	1715	PAR	C11-C21	2.50	1.57	1.52
58	XA	1715	PAR	C31-C21	2.43	1.56	1.53
58	QA	1696	PAR	O54-C54	2.22	1.49	1.44
58	QA	1696	PAR	C31-C21	2.15	1.56	1.53
58	QA	1696	PAR	C44-C54	2.12	1.57	1.53
58	QA	1696	PAR	C14-C24	2.05	1.56	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
58	QA	1696	PAR	C14-O54-C54	5.03	123.55	113.72
58	XA	1715	PAR	O33-C14-C24	4.87	116.06	108.08
58	XA	1715	PAR	C14-O54-C54	3.93	121.40	113.72
58	QA	1696	PAR	O33-C14-C24	3.32	113.51	108.08
58	QA	1696	PAR	O54-C54-C64	3.14	112.11	106.07
58	XA	1715	PAR	O54-C54-C64	2.96	111.75	106.07
58	QA	1696	PAR	C11-O51-C51	2.52	118.65	113.72
58	XA	1715	PAR	C11-O51-C51	2.30	118.21	113.72
58	QA	1696	PAR	C44-C34-C24	-2.17	107.39	110.99
58	QA	1696	PAR	C22-C32-C42	2.13	114.72	109.50

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
58	QA	1696	PAR	O43-C43-C53-O53
58	XA	1715	PAR	O43-C43-C53-O53
58	QA	1696	PAR	C33-C43-C53-O53
58	XA	1715	PAR	C33-C43-C53-O53
58	XA	1715	PAR	O51-C51-C61-O61
58	XA	1715	PAR	C41-C51-C61-O61
58	QA	1696	PAR	C52-C42-O11-C11
58	XA	1715	PAR	C52-C42-O11-C11
58	QA	1696	PAR	C44-C54-C64-N64
58	XA	1715	PAR	C43-C33-O33-C14

There are no ring outliers.

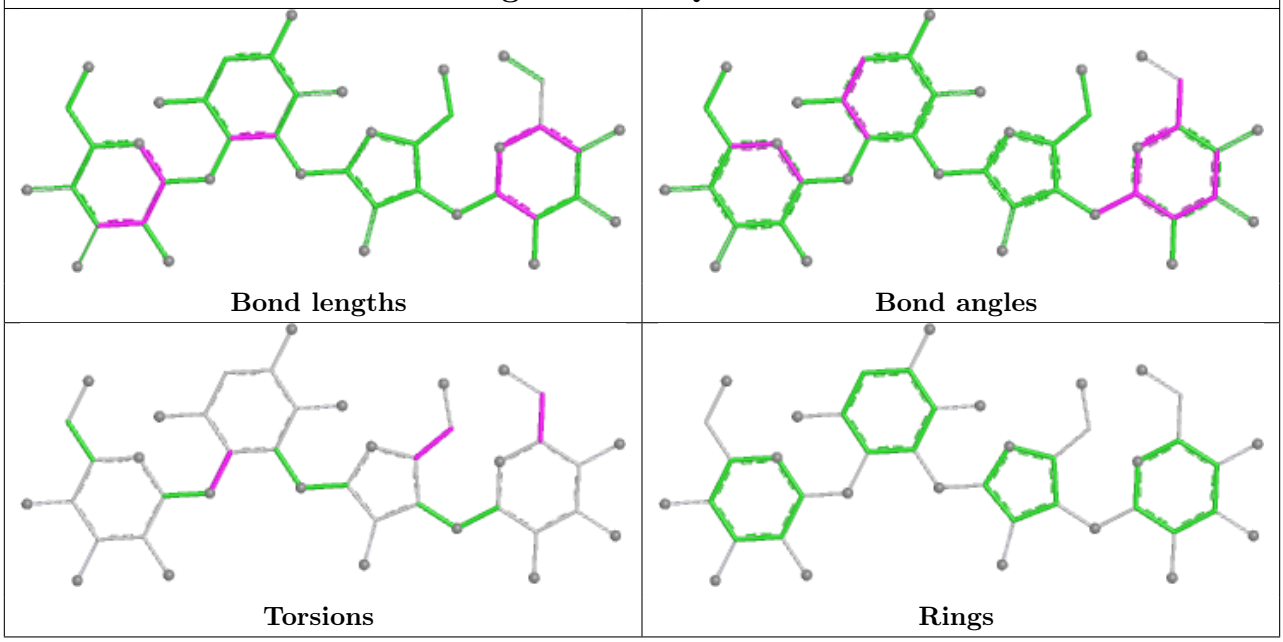
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
58	QA	1696	PAR	3	0
58	XA	1715	PAR	1	0

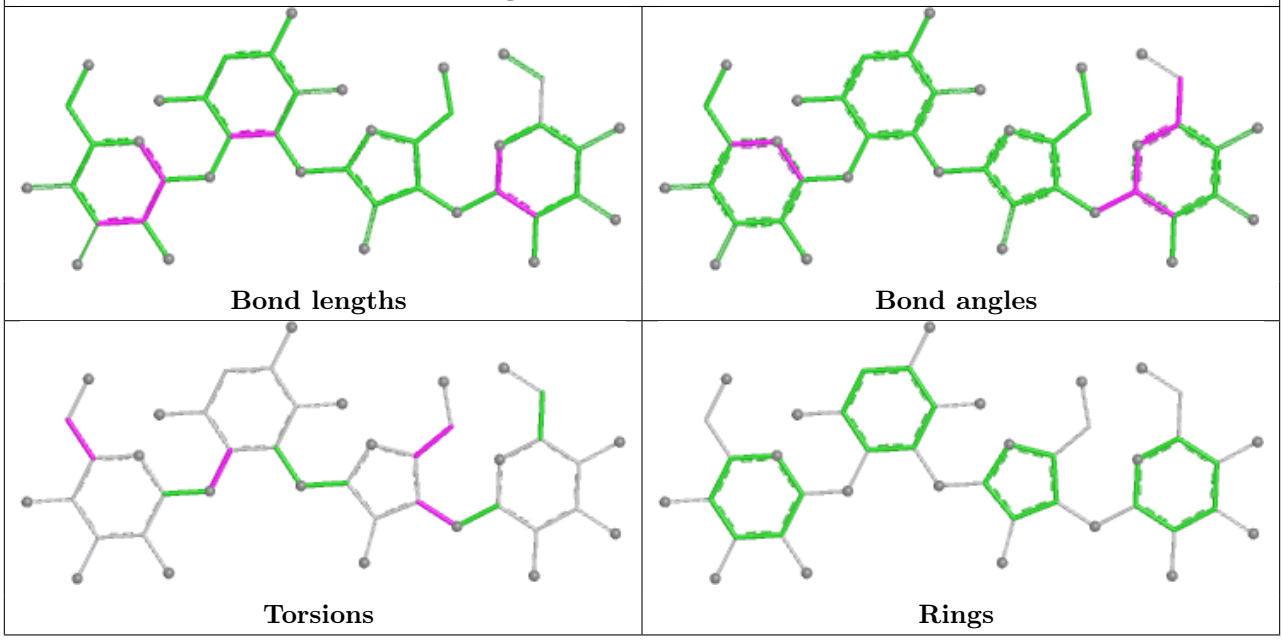
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand PAR QA 1696



Ligand PAR XA 1715



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	QA	1500/1522 (98%)	0.21	31 (2%) 63 37	37, 82, 166, 360	0
1	XA	1500/1522 (98%)	0.07	16 (1%) 78 52	27, 73, 170, 298	0
2	QB	237/256 (92%)	0.57	17 (7%) 21 14	55, 128, 200, 260	0
2	XB	237/256 (92%)	0.41	4 (1%) 69 41	51, 115, 181, 270	0
3	QC	205/239 (85%)	0.37	5 (2%) 59 34	64, 119, 182, 242	0
3	XC	205/239 (85%)	0.24	7 (3%) 48 27	45, 87, 134, 188	0
4	QD	208/209 (99%)	0.55	12 (5%) 29 17	39, 90, 141, 165	0
4	XD	208/209 (99%)	0.53	12 (5%) 29 17	34, 83, 139, 183	0
5	QE	151/162 (93%)	0.54	5 (3%) 49 28	49, 94, 145, 200	0
5	XE	151/162 (93%)	0.45	3 (1%) 65 38	39, 76, 128, 177	0
6	QF	101/101 (100%)	-0.16	0 100 100	34, 75, 114, 156	0
6	XF	101/101 (100%)	0.14	0 100 100	30, 77, 132, 236	0
7	QG	155/156 (99%)	0.36	5 (3%) 50 29	55, 111, 167, 291	0
7	XG	155/156 (99%)	0.36	4 (2%) 57 33	52, 101, 154, 205	0
8	QH	138/138 (100%)	0.52	5 (3%) 46 26	57, 93, 145, 158	0
8	XH	138/138 (100%)	0.41	3 (2%) 62 36	44, 83, 126, 144	0
9	QI	127/128 (99%)	1.07	17 (13%) 7 6	62, 122, 183, 230	0
9	XI	127/128 (99%)	1.04	17 (13%) 7 6	39, 114, 166, 213	0
10	QJ	99/105 (94%)	1.12	14 (14%) 6 6	56, 133, 221, 251	0
10	XJ	99/105 (94%)	0.63	3 (3%) 52 30	43, 113, 174, 209	0
11	QK	119/129 (92%)	0.39	5 (4%) 40 22	46, 83, 150, 237	0
11	XK	119/129 (92%)	0.16	4 (3%) 48 27	33, 76, 147, 227	0
12	QL	125/132 (94%)	1.06	20 (16%) 5 5	42, 80, 142, 248	0
12	XL	125/132 (94%)	0.80	20 (16%) 5 5	32, 68, 126, 266	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
13	QM	121/126 (96%)	0.92	13 (10%) 11 9	45, 118, 179, 283	0
13	XM	121/126 (96%)	0.68	10 (8%) 17 12	41, 99, 165, 284	0
14	QN	60/61 (98%)	1.03	6 (10%) 12 10	52, 104, 144, 159	0
14	XN	60/61 (98%)	0.70	6 (10%) 12 10	37, 76, 109, 126	0
15	QO	88/89 (98%)	0.21	0 100 100	36, 84, 136, 178	0
15	XO	88/89 (98%)	0.35	0 100 100	41, 76, 120, 130	0
16	QP	84/88 (95%)	0.82	5 (5%) 27 17	43, 78, 129, 174	0
16	XP	84/88 (95%)	0.90	7 (8%) 17 12	53, 89, 138, 202	0
17	QQ	100/105 (95%)	0.52	2 (2%) 65 38	51, 90, 130, 190	0
17	XQ	100/105 (95%)	0.53	5 (5%) 34 19	48, 85, 127, 173	0
18	QR	70/88 (79%)	0.00	0 100 100	43, 82, 138, 172	0
18	XR	70/88 (79%)	0.03	0 100 100	35, 79, 124, 157	0
19	QS	84/93 (90%)	0.68	4 (4%) 35 20	70, 127, 178, 246	0
19	XS	84/93 (90%)	0.45	3 (3%) 46 26	63, 106, 161, 228	0
20	QT	99/106 (93%)	0.96	12 (12%) 8 7	56, 98, 150, 177	0
20	XT	99/106 (93%)	1.14	16 (16%) 4 5	61, 110, 169, 201	0
21	QU	25/27 (92%)	1.49	6 (24%) 2 2	62, 110, 148, 166	0
21	XU	25/27 (92%)	1.08	4 (16%) 5 5	54, 90, 124, 164	0
22	QV	77/77 (100%)	0.25	1 (1%) 75 48	41, 95, 162, 189	0
22	XV	77/77 (100%)	0.20	1 (1%) 75 48	34, 87, 125, 169	0
23	QX	11/25 (44%)	1.52	3 (27%) 1 2	55, 79, 176, 213	0
23	XX	11/25 (44%)	1.65	4 (36%) 1 1	41, 59, 161, 175	0
24	QY	15/17 (88%)	0.18	0 100 100	62, 89, 153, 156	0
24	XY	16/17 (94%)	0.24	0 100 100	57, 85, 152, 154	0
25	RA	2882/2915 (98%)	0.02	37 (1%) 75 48	24, 60, 218, 384	0
25	YA	2882/2915 (98%)	-0.06	32 (1%) 78 52	19, 51, 215, 471	0
26	RB	120/122 (98%)	0.35	2 (1%) 69 41	69, 96, 130, 160	0
26	YB	120/122 (98%)	0.03	2 (1%) 69 41	50, 73, 106, 142	0
27	RD	272/276 (98%)	0.30	11 (4%) 42 24	22, 49, 98, 171	0
27	YD	272/276 (98%)	0.33	9 (3%) 49 28	11, 45, 84, 223	0
28	RE	205/206 (99%)	0.70	14 (6%) 23 15	22, 70, 140, 321	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
28	YE	205/206 (99%)	0.61	13 (6%) 26 16	22, 68, 139, 246	0
29	RF	202/210 (96%)	0.06	1 (0%) 87 66	20, 71, 132, 203	0
29	YF	202/210 (96%)	0.27	2 (0%) 79 54	14, 63, 129, 185	0
30	RG	181/182 (99%)	0.61	8 (4%) 39 22	64, 114, 167, 218	0
30	YG	181/182 (99%)	0.43	7 (3%) 43 24	50, 83, 136, 272	0
31	RH	170/180 (94%)	0.96	16 (9%) 14 10	77, 169, 247, 338	0
31	YH	170/180 (94%)	0.74	11 (6%) 25 16	36, 90, 147, 208	0
32	RI	146/148 (98%)	0.70	13 (8%) 15 11	37, 104, 161, 269	0
32	YI	146/148 (98%)	0.30	5 (3%) 48 27	26, 102, 156, 181	0
33	RN	138/140 (98%)	0.51	2 (1%) 73 46	38, 75, 133, 194	0
33	YN	138/140 (98%)	0.44	6 (4%) 40 22	36, 70, 124, 180	0
34	RO	122/122 (100%)	0.30	4 (3%) 49 28	28, 65, 109, 146	0
34	YO	122/122 (100%)	-0.00	1 (0%) 82 58	17, 51, 86, 108	0
35	RP	150/150 (100%)	0.41	7 (4%) 36 20	19, 79, 151, 264	0
35	YP	150/150 (100%)	0.51	8 (5%) 32 19	15, 68, 135, 202	0
36	RQ	141/141 (100%)	1.17	24 (17%) 4 4	43, 88, 145, 228	0
36	YQ	141/141 (100%)	0.88	21 (14%) 5 5	18, 63, 121, 227	0
37	RR	118/118 (100%)	0.26	2 (1%) 69 41	26, 57, 91, 135	0
37	YR	118/118 (100%)	0.45	1 (0%) 82 58	34, 65, 112, 156	0
38	RS	111/112 (99%)	0.83	10 (9%) 15 11	60, 96, 158, 199	0
38	YS	111/112 (99%)	0.89	11 (9%) 13 10	36, 84, 137, 172	0
39	RT	137/146 (93%)	0.51	8 (5%) 29 17	39, 74, 168, 287	0
39	YT	137/146 (93%)	0.35	6 (4%) 39 22	29, 71, 152, 217	0
40	RU	117/118 (99%)	0.34	1 (0%) 81 56	23, 68, 138, 196	0
40	YU	117/118 (99%)	0.35	5 (4%) 40 22	23, 57, 136, 223	0
41	RV	101/101 (100%)	0.36	2 (1%) 65 38	37, 89, 166, 286	0
41	YV	101/101 (100%)	0.52	8 (7%) 18 13	31, 84, 143, 275	0
42	RW	113/113 (100%)	0.33	4 (3%) 47 27	22, 52, 109, 206	0
42	YW	113/113 (100%)	0.38	4 (3%) 47 27	24, 53, 119, 205	0
43	RX	92/96 (95%)	0.43	4 (4%) 40 22	22, 65, 110, 134	0
43	YX	92/96 (95%)	0.39	2 (2%) 62 36	15, 53, 95, 119	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
44	RY	102/110 (92%)	1.21	20 (19%) 3 3	48, 95, 183, 309	0
44	YY	102/110 (92%)	0.76	12 (11%) 9 8	30, 79, 135, 264	0
45	RZ	183/206 (88%)	0.90	14 (7%) 19 13	65, 123, 209, 310	0
45	YZ	183/206 (88%)	0.76	12 (6%) 24 15	49, 101, 195, 273	0
46	R0	82/85 (96%)	0.60	3 (3%) 45 25	33, 71, 100, 128	0
46	Y0	82/85 (96%)	0.45	3 (3%) 45 25	29, 55, 90, 119	0
47	R1	97/98 (98%)	0.62	8 (8%) 17 12	25, 62, 140, 257	0
47	Y1	97/98 (98%)	0.65	7 (7%) 21 14	17, 56, 154, 232	0
48	R2	69/72 (95%)	0.47	2 (2%) 53 30	42, 81, 158, 197	0
48	Y2	69/72 (95%)	0.40	1 (1%) 73 46	30, 64, 139, 181	0
49	R3	59/60 (98%)	0.39	1 (1%) 69 41	45, 86, 134, 177	0
49	Y3	59/60 (98%)	0.35	1 (1%) 69 41	27, 60, 128, 185	0
50	R4	71/71 (100%)	1.36	14 (19%) 3 3	110, 208, 323, 389	0
50	Y4	71/71 (100%)	0.75	2 (2%) 55 31	66, 140, 233, 334	0
51	R5	59/60 (98%)	0.78	5 (8%) 16 12	23, 64, 219, 249	0
51	Y5	58/60 (96%)	0.80	7 (12%) 8 7	27, 71, 208, 261	0
52	R6	49/54 (90%)	1.23	7 (14%) 6 6	110, 164, 240, 272	0
52	Y6	49/54 (90%)	1.10	7 (14%) 6 6	95, 147, 224, 286	0
53	R7	49/49 (100%)	0.42	4 (8%) 17 12	24, 44, 95, 172	0
53	Y7	49/49 (100%)	0.38	3 (6%) 27 16	18, 37, 95, 166	0
54	R8	64/65 (98%)	1.21	11 (17%) 4 4	26, 68, 121, 209	0
54	Y8	64/65 (98%)	1.14	11 (17%) 4 4	27, 58, 128, 223	0
55	R9	37/37 (100%)	2.80	25 (67%) 0 0	96, 171, 236, 316	0
55	Y9	37/37 (100%)	2.41	21 (56%) 0 0	99, 151, 210, 312	0
56	Z5	2/3 (66%)	0.19	0 100 100	35, 35, 35, 59	0
56	Z6	2/3 (66%)	0.70	0 100 100	32, 32, 32, 42	0
All	All	20880/21490 (97%)	0.34	862 (4%) 41 23	11, 75, 176, 471	0

All (862) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
11	QK	128	ALA	8.0
23	XX	25	A	7.3

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Mol	Chain	Res	Type	RSRZ
53	R7	48	LYS	6.6
53	Y7	48	LYS	6.2
36	YQ	80	GLU	6.1
23	XX	15	A	5.8
11	QK	127	LYS	5.8
55	R9	11	CYS	5.8
54	Y8	64	TYR	5.8
14	XN	2	ALA	5.8
7	XG	5	ARG	5.8
47	Y1	98	LEU	5.7
23	QX	25	A	5.7
10	QJ	58	ASP	5.7
55	R9	14	CYS	5.7
12	QL	19	ARG	5.6
36	RQ	80	GLU	5.6
55	Y9	26	ILE	5.5
25	RA	34	C	5.5
55	R9	15	LYS	5.5
44	RY	29	GLU	5.4
13	XM	94	ARG	5.3
13	QM	122	LYS	5.2
9	QI	127	LYS	5.1
9	QI	115	GLY	5.1
53	Y7	49	ARG	5.1
31	YH	155	SER	5.1
53	R7	49	ARG	5.0
33	RN	68	GLU	5.0
55	Y9	28	GLU	4.9
9	XI	12	GLU	4.8
22	QV	0	C	4.7
32	RI	16	GLY	4.7
32	RI	15	VAL	4.7
20	XT	9	ASN	4.6
13	XM	122	LYS	4.6
11	XK	127	LYS	4.6
9	XI	110	GLU	4.6
21	QU	2	GLY	4.6
55	R9	16	VAL	4.5
1	QA	82	U	4.5
44	RY	62	GLU	4.5
13	QM	102	ARG	4.5
45	YZ	167	PRO	4.5

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Mol	Chain	Res	Type	RSRZ
1	XA	5	U	4.5
16	XP	68	ASP	4.4
44	YY	2	ARG	4.4
11	QK	129	SER	4.4
20	XT	70	SER	4.4
12	QL	64	TYR	4.4
55	R9	37	GLY	4.4
31	RH	169	VAL	4.4
35	YP	16	ARG	4.3
27	YD	235	GLY	4.3
22	XV	0	C	4.3
55	Y9	8	LYS	4.2
42	RW	22	ASP	4.2
55	Y9	12	ASP	4.2
50	R4	69	LYS	4.2
25	RA	6	A	4.1
54	R8	48	PHE	4.1
44	RY	43	ASN	4.1
36	RQ	87	LYS	4.1
54	R8	50	LEU	4.1
9	QI	30	GLY	4.1
36	RQ	88	GLY	4.1
10	QJ	45	ARG	4.1
10	XJ	59	SER	4.1
55	R9	36	GLN	4.1
4	QD	80	GLU	4.1
32	YI	15	VAL	4.1
8	QH	1	MET	4.1
25	YA	2188	C	4.0
20	XT	16	HIS	4.0
47	Y1	95	LEU	4.0
55	Y9	32	HIS	4.0
9	QI	128	ARG	4.0
11	XK	129	SER	3.9
25	YA	34	C	3.9
31	RH	155	SER	3.9
44	RY	58	GLY	3.9
54	R8	64	TYR	3.9
17	QQ	24	GLU	3.9
16	XP	48	TRP	3.9
55	R9	23	VAL	3.9
13	QM	10	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
9	QI	105	ASP	3.8
14	QN	46	GLU	3.8
36	RQ	91	GLU	3.8
1	XA	1531	A	3.8
31	RH	43	VAL	3.8
12	QL	28	LYS	3.8
39	YT	104	ASN	3.8
21	XU	10	ARG	3.8
44	YY	91	GLU	3.8
55	R9	26	ILE	3.8
31	RH	170	ARG	3.8
55	Y9	2	LYS	3.8
36	RQ	83	MET	3.8
50	R4	51	ASP	3.8
55	R9	2	LYS	3.8
4	XD	163	GLU	3.7
50	R4	68	ARG	3.7
36	RQ	20	ALA	3.7
12	XL	27	LEU	3.7
10	QJ	59	SER	3.7
35	RP	16	ARG	3.7
12	XL	7	ILE	3.7
41	RV	36	PRO	3.7
20	QT	12	ALA	3.7
36	RQ	33	GLY	3.7
44	RY	49	VAL	3.7
20	QT	22	ARG	3.7
12	QL	16	GLU	3.7
50	R4	49	PHE	3.7
12	XL	28	LYS	3.7
12	QL	23	LYS	3.6
39	YT	106	SER	3.6
25	YA	1176	G	3.6
46	R0	3	HIS	3.6
14	QN	35	ARG	3.6
27	RD	28	GLU	3.6
12	QL	129	ALA	3.6
28	RE	63	LEU	3.6
9	XI	106	ALA	3.6
33	YN	68	GLU	3.6
26	YB	1	U	3.6
35	RP	61	ARG	3.6

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Mol	Chain	Res	Type	RSRZ
4	XD	22	LYS	3.5
10	QJ	46	ARG	3.5
1	QA	1531	A	3.5
28	RE	21	VAL	3.5
25	RA	2145	C	3.5
12	QL	127	GLU	3.5
28	RE	204	ALA	3.5
36	RQ	21	THR	3.5
45	YZ	166	SER	3.5
52	Y6	23	THR	3.5
36	RQ	89	ASN	3.5
35	YP	15	ARG	3.5
41	YV	36	PRO	3.5
28	YE	200	GLU	3.5
41	YV	48	GLY	3.5
23	XX	24	A	3.5
4	XD	50	ARG	3.5
39	RT	115	ARG	3.5
42	RW	92	ARG	3.5
4	QD	32	ALA	3.4
12	QL	128	ALA	3.4
9	XI	126	SER	3.4
25	YA	2062	A	3.4
27	RD	44	ASN	3.4
55	R9	32	HIS	3.4
20	XT	19	SER	3.4
36	YQ	21	THR	3.4
23	QX	15	A	3.4
9	QI	106	ALA	3.4
12	QL	27	LEU	3.4
28	RE	159	HIS	3.4
54	R8	61	LEU	3.4
25	YA	2165	G	3.4
44	RY	63	LYS	3.4
39	YT	11	GLU	3.4
13	QM	6	GLY	3.4
4	QD	73	ARG	3.4
10	QJ	47	PHE	3.4
9	XI	115	GLY	3.4
9	XI	64	THR	3.4
27	RD	26	LYS	3.4
55	Y9	1	MET	3.4

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Mol	Chain	Res	Type	RSRZ
25	RA	1026	U	3.4
25	RA	2062	A	3.4
40	YU	90	VAL	3.4
52	Y6	5	VAL	3.4
12	QL	91	LYS	3.4
12	XL	64	TYR	3.4
17	XQ	101	ARG	3.4
47	R1	92	LYS	3.4
36	YQ	79	LEU	3.3
5	QE	94	ALA	3.3
55	Y9	29	ASN	3.3
36	YQ	10	ARG	3.3
36	YQ	87	LYS	3.3
32	RI	1	MET	3.3
36	RQ	18	LYS	3.3
36	YQ	91	GLU	3.3
4	QD	31	CYS	3.3
7	XG	8	GLU	3.3
44	RY	46	LYS	3.3
7	QG	5	ARG	3.3
27	RD	33	LEU	3.3
9	QI	110	GLU	3.3
12	QL	21	LYS	3.3
21	XU	26	LYS	3.3
38	RS	11	LYS	3.3
36	YQ	140	ALA	3.3
9	QI	126	SER	3.3
36	RQ	1	MET	3.3
5	XE	81	GLU	3.3
54	Y8	40	GLU	3.3
46	R0	2	ALA	3.2
55	R9	25	VAL	3.2
9	XI	75	ASP	3.2
25	YA	2189	U	3.2
26	RB	88	C	3.2
27	YD	40	THR	3.2
10	QJ	62	HIS	3.2
51	Y5	60	VAL	3.2
25	YA	2804	C	3.2
9	QI	27	THR	3.2
31	YH	58	GLU	3.2
51	R5	2	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
54	Y8	65	GLU	3.2
44	RY	103	GLY	3.2
55	Y9	34	GLN	3.2
8	XH	1	MET	3.2
12	XL	19	ARG	3.1
4	XD	26	CYS	3.1
10	QJ	64	GLU	3.1
40	RU	90	VAL	3.1
45	YZ	155	LEU	3.1
35	RP	15	ARG	3.1
12	XL	17	LYS	3.1
43	RX	38	GLU	3.1
20	XT	15	ARG	3.1
55	Y9	22	ARG	3.1
13	QM	101	GLN	3.1
41	YV	37	VAL	3.1
12	XL	16	GLU	3.1
25	YA	508	G	3.1
2	QB	79	ASP	3.1
52	R6	7	ILE	3.1
13	XM	100	GLY	3.1
55	Y9	21	GLY	3.1
1	QA	532	A	3.1
25	YA	2179	C	3.1
37	YR	102	GLU	3.1
20	XT	18	GLN	3.1
44	RY	102	CYS	3.1
28	RE	7	VAL	3.0
9	XI	119	ALA	3.0
52	R6	42	TRP	3.0
3	QC	161	GLU	3.0
16	QP	54	GLU	3.0
25	RA	614	U	3.0
30	YG	26	GLN	3.0
44	RY	52	SER	3.0
29	RF	131	GLY	3.0
31	RH	115	VAL	3.0
4	QD	50	ARG	3.0
30	YG	167	GLU	3.0
25	RA	2211	G	3.0
7	XG	81	GLY	3.0
25	YA	887	A	3.0

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Mol	Chain	Res	Type	RSRZ
4	XD	32	ALA	3.0
21	XU	2	GLY	3.0
20	QT	25	ARG	3.0
30	RG	91	ARG	3.0
1	QA	1181	G	3.0
25	RA	2121	G	3.0
54	Y8	48	PHE	3.0
11	XK	128	ALA	3.0
28	YE	204	ALA	3.0
3	XC	190	ARG	2.9
23	QX	24	A	2.9
1	XA	1453	G	2.9
25	YA	11	G	2.9
9	QI	117	HIS	2.9
5	XE	83	GLU	2.9
12	XL	5	PRO	2.9
44	YY	64	GLU	2.9
1	XA	1281	U	2.9
9	XI	8	GLY	2.9
47	R1	96	LYS	2.9
55	Y9	31	LYS	2.9
55	R9	17	ILE	2.9
52	Y6	42	TRP	2.9
31	RH	145	ALA	2.9
30	YG	80	PHE	2.9
38	YS	41	ASP	2.9
20	XT	69	GLY	2.9
20	QT	74	LYS	2.9
44	RY	33	LYS	2.9
8	QH	86	ILE	2.9
4	XD	31	CYS	2.9
12	QL	32	PHE	2.9
2	XB	12	GLU	2.9
1	QA	1348	U	2.9
1	XA	631	G	2.9
25	RA	12	U	2.9
25	RA	2164	C	2.9
7	QG	81	GLY	2.9
12	QL	17	LYS	2.9
21	QU	4	GLY	2.9
55	R9	31	LYS	2.9
55	Y9	15	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
49	Y3	55	ARG	2.9
20	QT	21	LYS	2.9
36	RQ	19	GLY	2.8
32	RI	116	LEU	2.8
41	YV	45	THR	2.8
50	R4	71	ARG	2.8
13	XM	115	LYS	2.8
13	QM	8	GLU	2.8
33	RN	10	GLU	2.8
2	QB	41	ILE	2.8
1	QA	1183	A	2.8
12	XL	13	LYS	2.8
12	XL	21	LYS	2.8
44	YY	63	LYS	2.8
55	R9	30	PRO	2.8
20	QT	11	SER	2.8
1	QA	723	U	2.8
1	XA	630	G	2.8
4	QD	209	ARG	2.8
28	YE	152	LYS	2.8
38	YS	33	LYS	2.8
27	RD	64	ILE	2.8
27	YD	33	LEU	2.8
47	Y1	97	LEU	2.8
54	Y8	61	LEU	2.8
9	QI	64	THR	2.8
25	YA	1177	A	2.8
47	Y1	92	LYS	2.8
9	XI	102	LEU	2.8
2	QB	4	GLU	2.8
27	RD	55	GLY	2.8
30	RG	28	VAL	2.8
51	Y5	52	TYR	2.8
2	QB	231	GLU	2.8
17	XQ	24	GLU	2.8
32	RI	146	ALA	2.8
21	XU	15	ARG	2.7
49	R3	30	ARG	2.7
36	RQ	90	VAL	2.7
39	RT	1	MET	2.7
55	R9	34	GLN	2.7
10	QJ	67	THR	2.7

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Mol	Chain	Res	Type	RSRZ
20	XT	35	THR	2.7
25	YA	275	G	2.7
52	R6	26	ASN	2.7
47	R1	95	LEU	2.7
25	YA	1536	A	2.7
9	QI	75	ASP	2.7
21	QU	5	ASP	2.7
12	QL	37	CYS	2.7
9	XI	114	TYR	2.7
12	QL	26	ALA	2.7
28	YE	193	GLY	2.7
31	YH	5	GLY	2.7
35	YP	74	GLU	2.7
1	QA	980	C	2.7
20	QT	9	ASN	2.7
21	QU	10	ARG	2.7
50	R4	66	SER	2.7
19	XS	76	PRO	2.7
25	YA	2112	G	2.7
25	YA	2141	G	2.7
13	QM	100	GLY	2.7
31	RH	171	LEU	2.7
33	YN	57	ALA	2.7
30	YG	35	GLU	2.7
43	YX	65	ARG	2.7
55	R9	35	ARG	2.7
20	XT	14	LYS	2.7
52	Y6	8	LYS	2.7
1	XA	422	C	2.7
38	YS	64	GLU	2.7
40	YU	89	GLU	2.7
45	YZ	162	GLU	2.7
3	QC	135	LYS	2.7
25	RA	1847	A	2.7
36	RQ	59	ARG	2.7
14	QN	55	GLY	2.7
33	YN	49	GLY	2.7
44	RY	44	ILE	2.7
27	RD	35	LYS	2.7
50	R4	38	LYS	2.7
12	QL	62	SER	2.7
31	YH	169	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
42	YW	92	ARG	2.6
54	R8	53	PRO	2.6
16	QP	65	GLN	2.6
16	XP	47	ASP	2.6
25	RA	2170	A	2.6
55	R9	13	LYS	2.6
27	YD	237	GLU	2.6
1	QA	1196	U	2.6
25	YA	12	U	2.6
52	R6	5	VAL	2.6
35	RP	24	GLY	2.6
38	YS	2	ALA	2.6
55	Y9	36	GLN	2.6
48	R2	52	ASP	2.6
50	R4	27	THR	2.6
51	Y5	51	TYR	2.6
55	R9	28	GLU	2.6
12	XL	96	VAL	2.6
12	XL	89	ARG	2.6
25	YA	1026	U	2.6
42	YW	113	LYS	2.6
38	RS	2	ALA	2.6
54	R8	35	GLN	2.6
4	QD	23	GLY	2.6
38	RS	7	TYR	2.6
20	XT	23	ARG	2.6
30	YG	52	ILE	2.6
32	YI	79	ILE	2.6
36	YQ	20	ALA	2.6
3	XC	167	TRP	2.6
19	QS	80	TYR	2.6
50	R4	67	TYR	2.6
42	RW	67	ASP	2.6
13	XM	7	VAL	2.6
31	RH	52	VAL	2.6
39	RT	112	ARG	2.6
39	YT	111	ARG	2.6
45	RZ	20	ARG	2.6
44	RY	47	LYS	2.6
12	XL	128	ALA	2.6
28	YE	78	LEU	2.6
33	YN	138	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
55	R9	20	HIS	2.6
28	YE	107	THR	2.6
3	QC	167	TRP	2.6
3	XC	161	GLU	2.6
9	QI	12	GLU	2.6
38	RS	64	GLU	2.6
48	Y2	11	GLU	2.6
39	RT	106	SER	2.6
26	YB	88	C	2.6
13	XM	10	PRO	2.6
12	XL	95	GLY	2.6
42	RW	112	GLY	2.6
28	RE	76	ARG	2.5
31	YH	18	GLU	2.5
39	RT	46	GLU	2.5
39	YT	109	GLU	2.5
54	Y8	54	GLU	2.5
16	QP	68	ASP	2.5
45	RZ	87	ASP	2.5
28	RE	161	GLY	2.5
28	RE	193	GLY	2.5
25	RA	1092	C	2.5
52	R6	23	THR	2.5
55	R9	10	ILE	2.5
20	QT	23	ARG	2.5
55	R9	22	ARG	2.5
1	QA	1450	U	2.5
25	RA	1420	U	2.5
25	RA	1460	A	2.5
28	YE	63	LEU	2.5
20	XT	75	ASN	2.5
46	Y0	2	ALA	2.5
2	XB	133	LYS	2.5
12	XL	91	LYS	2.5
2	QB	130	ARG	2.5
55	Y9	20	HIS	2.5
41	RV	81	TYR	2.5
50	R4	37	SER	2.5
44	YY	49	VAL	2.5
25	RA	1176	G	2.5
25	YA	889	C	2.5
36	YQ	138	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
37	RR	69	ASP	2.5
14	XN	14	PRO	2.5
7	QG	82	GLY	2.5
27	YD	38	LYS	2.5
1	XA	1126	U	2.5
17	QQ	101	ARG	2.5
9	XI	41	VAL	2.5
2	QB	163	PHE	2.5
2	QB	160	ASP	2.5
21	QU	26	LYS	2.5
36	RQ	12	GLN	2.5
44	YY	46	LYS	2.5
45	YZ	120	ILE	2.5
19	XS	84	GLY	2.5
33	YN	36	GLY	2.5
2	XB	217	ARG	2.5
27	YD	176	ARG	2.5
53	R7	23	ARG	2.5
20	XT	73	HIS	2.5
55	Y9	25	VAL	2.5
25	RA	2146	C	2.5
25	YA	2108	C	2.5
1	QA	1202	G	2.5
25	RA	2116	G	2.5
25	YA	274	G	2.5
34	RO	9	GLU	2.5
55	R9	33	LYS	2.5
4	QD	124	GLY	2.5
12	QL	72	GLY	2.5
52	R6	6	ARG	2.5
55	R9	9	ARG	2.5
54	R8	23	VAL	2.5
9	QI	125	TYR	2.5
10	QJ	98	ILE	2.4
10	QJ	85	LEU	2.4
9	QI	66	ARG	2.4
36	YQ	1	MET	2.4
1	QA	81	G	2.4
38	YS	59	LYS	2.4
3	QC	166	GLU	2.4
16	XP	16	HIS	2.4
29	YF	56	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
45	RZ	55	HIS	2.4
32	RI	12	LEU	2.4
4	XD	76	ARG	2.4
36	YQ	24	GLY	2.4
12	XL	23	LYS	2.4
32	RI	18	VAL	2.4
36	RQ	22	LYS	2.4
38	YS	11	LYS	2.4
4	QD	26	CYS	2.4
3	XC	166	GLU	2.4
32	RI	41	GLU	2.4
9	XI	66	ARG	2.4
13	QM	76	ALA	2.4
37	RR	2	ARG	2.4
25	RA	1093	G	2.4
25	RA	2319	G	2.4
16	XP	53	VAL	2.4
38	YS	69	VAL	2.4
10	XJ	58	ASP	2.4
50	Y4	5	ILE	2.4
4	QD	157	LEU	2.4
28	YE	151	TYR	2.4
39	RT	3	ARG	2.4
48	R2	46	GLN	2.4
52	R6	22	ALA	2.4
13	QM	106	ASN	2.4
20	QT	26	ASN	2.4
36	RQ	15	GLY	2.4
45	RZ	46	LYS	2.4
31	RH	113	VAL	2.4
55	Y9	23	VAL	2.4
47	R1	98	LEU	2.4
5	QE	81	GLU	2.4
17	XQ	91	ARG	2.4
27	RD	30	GLU	2.4
47	Y1	26	ARG	2.4
20	QT	16	HIS	2.4
38	YS	37	ALA	2.4
25	RA	1058	G	2.4
25	RA	1059	G	2.4
44	RY	34	LYS	2.4
44	YY	81	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
4	XD	89	THR	2.4
44	YY	90	LEU	2.4
14	XN	8	GLU	2.4
30	RG	30	GLU	2.4
51	Y5	59	GLU	2.4
25	YA	2144	U	2.3
4	QD	2	GLY	2.3
55	R9	21	GLY	2.3
1	QA	934	C	2.3
25	RA	2474	C	2.3
26	RB	12	C	2.3
30	RG	52	ILE	2.3
2	QB	165	VAL	2.3
31	YH	19	VAL	2.3
35	YP	50	ARG	2.3
35	YP	65	ARG	2.3
53	Y7	1	MET	2.3
1	QA	631	G	2.3
1	XA	81	G	2.3
25	YA	2121	G	2.3
25	YA	2190	G	2.3
38	RS	33	LYS	2.3
54	Y8	35	GLN	2.3
47	R1	85	LEU	2.3
1	QA	1126	U	2.3
25	RA	161	U	2.3
36	RQ	7	MET	2.3
54	R8	46	ARG	2.3
4	XD	30	LYS	2.3
11	QK	11	LYS	2.3
19	XS	85	LYS	2.3
54	Y8	52	LYS	2.3
1	QA	1286	A	2.3
23	XX	23	A	2.3
10	XJ	64	GLU	2.3
13	XM	16	ASP	2.3
20	QT	18	GLN	2.3
36	YQ	112	GLU	2.3
45	YZ	84	GLU	2.3
44	YY	89	PHE	2.3
3	XC	2	GLY	2.3
12	QL	31	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
54	Y8	63	PRO	2.3
27	RD	27	THR	2.3
29	YF	48	THR	2.3
12	XL	15	ARG	2.3
31	YH	3	ARG	2.3
34	RO	1	MET	2.3
25	YA	530	G	2.3
28	RE	73	GLU	2.3
33	YN	50	ASP	2.3
39	RT	11	GLU	2.3
42	YW	66	GLU	2.3
51	R5	59	GLU	2.3
54	R8	54	GLU	2.3
25	RA	9	U	2.3
25	RA	2122	U	2.3
38	YS	27	SER	2.3
20	QT	73	HIS	2.3
3	XC	165	THR	2.3
19	QS	79	THR	2.3
36	YQ	89	ASN	2.3
16	QP	48	TRP	2.3
34	RO	18	LYS	2.3
36	RQ	10	ARG	2.3
4	XD	12	CYS	2.3
32	RI	10	GLU	2.3
36	RQ	78	PRO	2.3
36	YQ	135	ASP	2.3
19	QS	11	VAL	2.3
38	RS	60	GLY	2.3
3	QC	30	ARG	2.3
28	RE	37	ARG	2.3
31	YH	6	ARG	2.3
1	QA	306	G	2.3
51	Y5	58	LEU	2.3
32	RI	17	GLN	2.3
51	Y5	46	CYS	2.3
1	QA	1346	A	2.3
14	XN	46	GLU	2.3
32	YI	117	GLU	2.3
1	QA	88	C	2.3
9	QI	28	VAL	2.3
14	QN	15	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
28	YE	56	PRO	2.3
45	YZ	130	PRO	2.3
7	QG	32	ARG	2.3
28	YE	1	MET	2.3
31	YH	152	ARG	2.3
32	RI	20	ASP	2.3
55	Y9	17	ILE	2.2
52	Y6	22	ALA	2.2
45	YZ	52	SER	2.2
27	YD	30	GLU	2.2
45	RZ	2	GLU	2.2
1	QA	1453	G	2.2
5	QE	99	GLY	2.2
8	QH	55	GLY	2.2
28	RE	22	PRO	2.2
38	RS	98	VAL	2.2
25	RA	1068	G	2.2
25	RA	2123	G	2.2
44	RY	59	GLY	2.2
53	R7	41	ARG	2.2
10	QJ	65	LEU	2.2
17	XQ	13	ASP	2.2
46	R0	10	THR	2.2
38	RS	34	HIS	2.2
39	RT	104	ASN	2.2
36	RQ	136	ALA	2.2
1	XA	328	C	2.2
50	R4	1	MET	2.2
2	QB	12	GLU	2.2
5	QE	24	ARG	2.2
8	QH	104	ARG	2.2
11	XK	12	ARG	2.2
13	QM	7	VAL	2.2
16	XP	71	ARG	2.2
35	RP	74	GLU	2.2
40	YU	111	GLU	2.2
43	RX	65	ARG	2.2
43	YX	60	ARG	2.2
54	Y8	46	ARG	2.2
28	RE	78	LEU	2.2
31	RH	154	PRO	2.2
40	YU	97	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
2	QB	16	HIS	2.2
1	QA	1065	U	2.2
1	QA	1235	U	2.2
30	YG	74	LYS	2.2
44	YY	101	LYS	2.2
55	Y9	13	LYS	2.2
46	Y0	17	GLN	2.2
36	YQ	7	MET	2.2
1	QA	1190	G	2.2
1	XA	1182	G	2.2
2	QB	144	ARG	2.2
9	XI	128	ARG	2.2
20	XT	20	LEU	2.2
28	YE	171	GLU	2.2
32	YI	7	GLU	2.2
36	YQ	48	GLU	2.2
51	R5	45	VAL	2.2
51	R5	50	GLY	2.2
25	RA	2896	C	2.2
2	XB	160	ASP	2.2
27	YD	4	LYS	2.2
28	YE	18	ASP	2.2
20	XT	12	ALA	2.2
19	QS	15	LEU	2.2
35	RP	18	ARG	2.2
36	YQ	5	ARG	2.2
45	RZ	125	LEU	2.2
12	XL	11	VAL	2.2
4	XD	80	GLU	2.2
31	YH	154	PRO	2.2
32	RI	122	GLU	2.2
42	YW	91	GLY	2.2
43	RX	94	GLY	2.2
50	R4	70	GLY	2.2
9	XI	78	LYS	2.2
44	RY	88	LYS	2.2
7	QG	152	ALA	2.2
35	RP	27	HIS	2.2
39	YT	1	MET	2.2
20	XT	72	LEU	2.2
25	RA	71	A	2.2
25	RA	1022	G	2.2

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Mol	Chain	Res	Type	RSRZ
14	QN	41	ARG	2.2
38	YS	20	ARG	2.2
8	XH	93	VAL	2.2
41	YV	58	VAL	2.2
25	YA	277	C	2.2
25	YA	888	C	2.2
43	RX	15	GLU	2.2
45	RZ	11	GLU	2.2
13	XM	121	LYS	2.2
38	RS	5	THR	2.2
28	RE	60	ASN	2.2
27	RD	37	LEU	2.2
36	YQ	83	MET	2.2
45	YZ	113	ALA	2.1
44	RY	6	HIS	2.1
44	YY	99	CYS	2.1
46	Y0	3	HIS	2.1
1	QA	1281	U	2.1
3	XC	193	TYR	2.1
31	RH	157	TYR	2.1
31	YH	45	VAL	2.1
44	RY	45	VAL	2.1
50	R4	31	ILE	2.1
52	Y6	7	ILE	2.1
45	RZ	78	LYS	2.1
45	YZ	114	GLY	2.1
30	YG	30	GLU	2.1
32	YI	41	GLU	2.1
1	QA	873	A	2.1
25	RA	529	A	2.1
25	RA	1057	A	2.1
1	QA	570	G	2.1
38	RS	32	LEU	2.1
47	R1	97	LEU	2.1
47	Y1	80	LEU	2.1
11	QK	89	ALA	2.1
10	QJ	43	ARG	2.1
16	QP	29	ASP	2.1
40	YU	81	HIS	2.1
2	QB	136	VAL	2.1
9	XI	108	VAL	2.1
34	RO	21	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
35	YP	64	LYS	2.1
12	QL	63	GLY	2.1
41	YV	101	GLY	2.1
54	R8	65	GLU	2.1
8	XH	3	THR	2.1
14	QN	22	THR	2.1
25	YA	614	U	2.1
12	XL	129	ALA	2.1
47	Y1	21	ARG	2.1
2	QB	68	ILE	2.1
2	QB	40	HIS	2.1
8	QH	97	VAL	2.1
27	RD	34	VAL	2.1
35	YP	108	LYS	2.1
25	YA	529	A	2.1
1	QA	413	G	2.1
1	XA	108	G	2.1
25	YA	2805	G	2.1
25	RA	2179	C	2.1
36	YQ	51	ARG	2.1
4	XD	164	ALA	2.1
13	QM	2	ALA	2.1
7	XG	84	ASN	2.1
1	XA	1125	U	2.1
45	RZ	105	VAL	2.1
45	RZ	116	VAL	2.1
55	Y9	3	VAL	2.1
9	XI	30	GLY	2.1
12	QL	95	GLY	2.1
30	RG	62	LEU	2.1
31	RH	88	LEU	2.1
4	QD	29	PRO	2.1
45	RZ	159	PRO	2.1
10	QJ	61	GLU	2.1
30	RG	35	GLU	2.1
36	RQ	48	GLU	2.1
45	RZ	119	GLU	2.1
44	RY	50	ARG	2.1
54	Y8	57	ARG	2.1
1	XA	1451	A	2.1
41	YV	51	VAL	2.1
1	QA	1359	C	2.1

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Mol	Chain	Res	Type	RSRZ
25	RA	2188	C	2.1
45	YZ	121	HIS	2.1
1	QA	1178	G	2.1
1	XA	64	G	2.1
31	RH	100	GLY	2.1
51	Y5	17	ASP	2.1
36	RQ	112	GLU	2.1
50	R4	34	GLU	2.1
44	YY	86	ARG	2.1
10	QJ	4	ILE	2.1
14	XN	58	LYS	2.1
16	XP	35	LYS	2.1
14	XN	59	ALA	2.1
31	RH	144	VAL	2.1
55	Y9	7	VAL	2.1
5	XE	85	GLY	2.0
31	RH	163	TYR	2.0
51	R5	34	PRO	2.0
13	QM	32	GLU	2.0
21	QU	17	THR	2.0
47	R1	26	ARG	2.0
55	R9	12	ASP	2.0
20	XT	34	LYS	2.0
30	RG	137	GLU	2.0
32	RI	118	LYS	2.0
34	YO	9	GLU	2.0
41	YV	53	GLU	2.0
2	QB	155	LEU	2.0
9	QI	124	GLN	2.0
25	RA	2183	C	2.0
1	QA	1182	G	2.0
25	YA	1678	G	2.0
31	RH	108	GLY	2.0
44	RY	55	TYR	2.0
50	Y4	22	ILE	2.0
52	Y6	27	LYS	2.0
45	YZ	97	GLU	2.0
54	R8	40	GLU	2.0
2	QB	186	ALA	2.0
12	XL	26	ALA	2.0
27	YD	15	PHE	2.0
13	QM	70	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
17	XQ	98	LEU	2.0
28	RE	20	ALA	2.0
5	QE	90	VAL	2.0
36	YQ	90	VAL	2.0
13	XM	102	ARG	2.0
13	XM	110	ARG	2.0
35	YP	61	ARG	2.0
38	YS	71	ARG	2.0
45	RZ	4	ARG	2.0
47	R1	79	GLY	2.0
28	YE	159	HIS	2.0
25	YA	2803	C	2.0
2	QB	181	PHE	2.0
30	RG	156	ASP	2.0
36	RQ	31	ASP	2.0
1	QA	1212	U	2.0
1	XA	723	U	2.0
25	RA	2189	U	2.0
45	RZ	172	ALA	2.0
36	YQ	81	VAL	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
56	PPU	Z5	101	37/38	0.92	0.14	44,44,45,45	0
56	PPU	Z6	101	37/38	0.93	0.12	28,28,29,29	0

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3128	1/1	0.42	0.23	24,24,24,24	0
57	MG	YA	3185	1/1	0.46	0.33	77,77,77,77	0
57	MG	YA	3308	1/1	0.47	0.35	64,64,64,64	0
57	MG	RA	3281	1/1	0.51	0.19	57,57,57,57	0
57	MG	RB	203	1/1	0.60	0.32	58,58,58,58	0
57	MG	YA	3272	1/1	0.60	0.21	37,37,37,37	0
57	MG	RA	3266	1/1	0.60	0.21	22,22,22,22	0
57	MG	RA	3198	1/1	0.62	0.38	46,46,46,46	0
57	MG	YA	3203	1/1	0.62	0.21	40,40,40,40	0
57	MG	QY	101	1/1	0.63	0.20	51,51,51,51	0
57	MG	YA	3173	1/1	0.63	0.31	40,40,40,40	0
57	MG	XA	1713	1/1	0.63	0.36	57,57,57,57	0
57	MG	R0	101	1/1	0.65	0.20	31,31,31,31	0
57	MG	YA	3300	1/1	0.66	0.36	69,69,69,69	0
57	MG	XA	1685	1/1	0.66	0.16	51,51,51,51	0
57	MG	RA	3301	1/1	0.67	0.21	52,52,52,52	0
57	MG	QX	101	1/1	0.67	0.22	38,38,38,38	0
57	MG	YA	3220	1/1	0.68	0.28	60,60,60,60	0
57	MG	RD	301	1/1	0.68	0.20	51,51,51,51	0
57	MG	QA	1630	1/1	0.69	0.15	80,80,80,80	0
57	MG	YA	3074	1/1	0.69	0.43	32,32,32,32	0
57	MG	YA	3295	1/1	0.70	0.28	62,62,62,62	0
57	MG	YA	3238	1/1	0.70	0.17	34,34,34,34	0
57	MG	QA	1695	1/1	0.70	0.27	44,44,44,44	0
57	MG	XA	1707	1/1	0.71	0.18	40,40,40,40	0
57	MG	XA	1712	1/1	0.71	0.20	49,49,49,49	0
57	MG	YA	3213	1/1	0.71	0.31	29,29,29,29	0
57	MG	YA	3216	1/1	0.71	0.24	47,47,47,47	0
57	MG	XA	1618	1/1	0.71	0.33	37,37,37,37	0
57	MG	RA	3210	1/1	0.72	0.18	29,29,29,29	0
57	MG	RA	3171	1/1	0.72	0.21	57,57,57,57	0
57	MG	RA	3268	1/1	0.72	0.18	47,47,47,47	0
57	MG	YA	3206	1/1	0.72	0.20	55,55,55,55	0
57	MG	YA	3337	1/1	0.72	0.34	41,41,41,41	0
57	MG	YA	3150	1/1	0.73	0.18	42,42,42,42	0
57	MG	YA	3320	1/1	0.73	0.30	50,50,50,50	0
57	MG	RA	3158	1/1	0.73	0.22	52,52,52,52	0
57	MG	RA	3262	1/1	0.74	0.21	38,38,38,38	0
57	MG	QA	1621	1/1	0.74	0.19	50,50,50,50	0
57	MG	YA	3296	1/1	0.74	0.26	55,55,55,55	0
57	MG	QA	1628	1/1	0.74	0.16	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3162	1/1	0.74	0.24	46,46,46,46	0
57	MG	YA	3317	1/1	0.74	0.21	57,57,57,57	0
57	MG	RA	3252	1/1	0.74	0.18	35,35,35,35	0
57	MG	YA	3323	1/1	0.74	0.21	52,52,52,52	0
57	MG	R8	101	1/1	0.74	0.17	29,29,29,29	0
57	MG	Y0	101	1/1	0.74	0.15	16,16,16,16	0
57	MG	RB	205	1/1	0.75	0.23	41,41,41,41	0
57	MG	YA	3328	1/1	0.75	0.31	54,54,54,54	0
57	MG	XA	1714	1/1	0.75	0.14	26,26,26,26	0
57	MG	RA	3287	1/1	0.75	0.28	60,60,60,60	0
57	MG	YA	3031	1/1	0.76	0.43	15,15,15,15	0
57	MG	QA	1625	1/1	0.76	0.14	59,59,59,59	0
57	MG	RA	3277	1/1	0.76	0.22	51,51,51,51	0
57	MG	YA	3174	1/1	0.76	0.23	50,50,50,50	0
57	MG	YA	3289	1/1	0.76	0.33	62,62,62,62	0
57	MG	RA	3290	1/1	0.77	0.11	62,62,62,62	0
57	MG	XA	1711	1/1	0.77	0.18	47,47,47,47	0
57	MG	RA	3244	1/1	0.77	0.51	60,60,60,60	0
57	MG	RA	3255	1/1	0.77	0.72	76,76,76,76	0
57	MG	YA	3285	1/1	0.77	0.28	31,31,31,31	0
57	MG	YA	3316	1/1	0.77	0.12	28,28,28,28	0
57	MG	XA	1658	1/1	0.78	0.10	61,61,61,61	0
57	MG	YA	3120	1/1	0.78	0.14	12,12,12,12	0
57	MG	YA	3302	1/1	0.78	0.25	41,41,41,41	0
57	MG	RA	3306	1/1	0.78	0.27	42,42,42,42	0
57	MG	YA	3311	1/1	0.78	0.21	47,47,47,47	0
57	MG	RA	3251	1/1	0.78	0.20	23,23,23,23	0
57	MG	YA	3221	1/1	0.78	0.21	28,28,28,28	0
57	MG	RA	3265	1/1	0.78	0.48	72,72,72,72	0
57	MG	RA	3010	1/1	0.78	0.15	37,37,37,37	0
57	MG	QA	1635	1/1	0.78	0.18	44,44,44,44	0
57	MG	RA	3291	1/1	0.78	0.23	30,30,30,30	0
57	MG	YB	202	1/1	0.78	0.16	49,49,49,49	0
57	MG	RA	3269	1/1	0.78	0.15	35,35,35,35	0
57	MG	YA	3307	1/1	0.79	0.20	40,40,40,40	0
57	MG	RA	3149	1/1	0.79	0.18	21,21,21,21	0
57	MG	YA	3280	1/1	0.79	0.42	67,67,67,67	0
57	MG	RA	3264	1/1	0.79	0.21	50,50,50,50	0
57	MG	YA	3218	1/1	0.79	0.15	40,40,40,40	0
57	MG	YA	3292	1/1	0.79	0.23	47,47,47,47	0
57	MG	RA	3293	1/1	0.79	0.20	42,42,42,42	0
57	MG	YA	3156	1/1	0.79	0.20	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3270	1/1	0.79	0.34	49,49,49,49	0
57	MG	YA	3267	1/1	0.79	0.19	41,41,41,41	0
57	MG	YA	3305	1/1	0.79	0.19	42,42,42,42	0
57	MG	YA	3322	1/1	0.80	0.25	44,44,44,44	0
57	MG	YA	3297	1/1	0.80	0.25	53,53,53,53	0
57	MG	YA	3116	1/1	0.80	0.15	13,13,13,13	0
57	MG	RB	202	1/1	0.80	0.18	59,59,59,59	0
57	MG	YB	201	1/1	0.80	0.31	47,47,47,47	0
57	MG	QA	1690	1/1	0.80	0.16	29,29,29,29	0
57	MG	YA	3187	1/1	0.80	0.21	37,37,37,37	0
57	MG	YA	3167	1/1	0.81	0.20	57,57,57,57	0
57	MG	YA	3293	1/1	0.81	0.21	40,40,40,40	0
57	MG	RA	3072	1/1	0.81	0.28	38,38,38,38	0
57	MG	RA	3246	1/1	0.81	0.34	48,48,48,48	0
57	MG	YA	3180	1/1	0.81	0.15	42,42,42,42	0
57	MG	RA	3250	1/1	0.81	0.18	25,25,25,25	0
57	MG	RA	3091	1/1	0.81	0.13	30,30,30,30	0
57	MG	QV	101	1/1	0.81	0.14	21,21,21,21	0
57	MG	QA	1682	1/1	0.81	0.09	20,20,20,20	0
57	MG	RA	3161	1/1	0.81	0.24	36,36,36,36	0
57	MG	XV	101	1/1	0.81	0.19	25,25,25,25	0
57	MG	QA	1619	1/1	0.81	0.12	40,40,40,40	0
57	MG	RA	3187	1/1	0.81	0.24	56,56,56,56	0
57	MG	QA	1667	1/1	0.81	0.26	44,44,44,44	0
57	MG	YA	3321	1/1	0.81	0.22	36,36,36,36	0
57	MG	RA	3202	1/1	0.81	0.29	50,50,50,50	0
57	MG	RA	3068	1/1	0.81	0.16	28,28,28,28	0
57	MG	YA	3148	1/1	0.81	0.16	55,55,55,55	0
57	MG	RA	3211	1/1	0.81	0.17	56,56,56,56	0
57	MG	YA	3284	1/1	0.81	0.16	39,39,39,39	0
57	MG	RA	3214	1/1	0.81	0.11	29,29,29,29	0
57	MG	XA	1656	1/1	0.81	0.20	52,52,52,52	0
57	MG	YA	3277	1/1	0.82	0.20	30,30,30,30	0
57	MG	YA	3278	1/1	0.82	0.18	32,32,32,32	0
57	MG	RA	3275	1/1	0.82	0.29	50,50,50,50	0
57	MG	RA	3013	1/1	0.82	0.26	25,25,25,25	0
57	MG	RA	3194	1/1	0.82	0.14	59,59,59,59	0
57	MG	RA	3014	1/1	0.82	0.37	29,29,29,29	0
57	MG	QA	1692	1/1	0.82	0.30	73,73,73,73	0
57	MG	QA	1679	1/1	0.83	0.26	45,45,45,45	0
57	MG	YA	3192	1/1	0.83	0.23	12,12,12,12	0
57	MG	RA	3307	1/1	0.83	0.61	75,75,75,75	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3267	1/1	0.83	0.28	57,57,57,57	0
57	MG	YA	3288	1/1	0.83	0.23	38,38,38,38	0
57	MG	RA	3286	1/1	0.83	0.13	66,66,66,66	0
57	MG	RA	3257	1/1	0.83	0.23	58,58,58,58	0
57	MG	RA	3175	1/1	0.83	0.15	24,24,24,24	0
57	MG	RA	3052	1/1	0.83	0.14	1,1,1,1	0
57	MG	XF	201	1/1	0.83	0.13	49,49,49,49	0
57	MG	YA	3335	1/1	0.83	0.15	19,19,19,19	0
57	MG	RA	3274	1/1	0.83	0.22	52,52,52,52	0
57	MG	YA	3029	1/1	0.83	0.31	6,6,6,6	0
57	MG	RA	3295	1/1	0.83	0.31	51,51,51,51	0
57	MG	RA	3208	1/1	0.83	0.08	32,32,32,32	0
57	MG	YA	3065	1/1	0.84	0.25	8,8,8,8	0
57	MG	QA	1654	1/1	0.84	0.14	46,46,46,46	0
57	MG	RA	3009	1/1	0.84	0.08	22,22,22,22	0
57	MG	YA	3119	1/1	0.84	0.13	15,15,15,15	0
57	MG	RA	3261	1/1	0.84	0.39	135,135,135,135	0
57	MG	YA	3274	1/1	0.84	0.27	45,45,45,45	0
57	MG	XA	1666	1/1	0.84	0.15	48,48,48,48	0
57	MG	RA	3237	1/1	0.84	0.14	32,32,32,32	0
57	MG	XA	1686	1/1	0.84	0.15	53,53,53,53	0
57	MG	XX	101	1/1	0.84	0.20	47,47,47,47	0
57	MG	XA	1695	1/1	0.84	0.16	43,43,43,43	0
57	MG	RA	3083	1/1	0.84	0.25	2,2,2,2	0
57	MG	QA	1620	1/1	0.85	0.21	40,40,40,40	0
57	MG	RA	3221	1/1	0.85	0.22	36,36,36,36	0
57	MG	RA	3303	1/1	0.85	0.27	50,50,50,50	0
57	MG	RA	3253	1/1	0.85	0.32	77,77,77,77	0
57	MG	YA	3137	1/1	0.85	0.17	12,12,12,12	0
57	MG	RA	3285	1/1	0.85	0.17	51,51,51,51	0
57	MG	YA	3209	1/1	0.85	0.17	34,34,34,34	0
57	MG	RA	3207	1/1	0.85	0.19	43,43,43,43	0
57	MG	RA	3062	1/1	0.85	0.16	51,51,51,51	0
57	MG	RA	3023	1/1	0.85	0.28	2,2,2,2	0
57	MG	YA	3290	1/1	0.85	0.28	41,41,41,41	0
57	MG	YA	3326	1/1	0.85	0.33	69,69,69,69	0
57	MG	YA	3045	1/1	0.85	0.25	5,5,5,5	0
57	MG	RA	3138	1/1	0.85	0.26	62,62,62,62	0
57	MG	YA	3233	1/1	0.85	0.18	32,32,32,32	0
57	MG	RA	3263	1/1	0.85	0.24	49,49,49,49	0
57	MG	YA	3257	1/1	0.85	0.28	9,9,9,9	0
57	MG	YA	3263	1/1	0.85	0.17	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3301	1/1	0.86	0.19	52,52,52,52	0
57	MG	RA	3152	1/1	0.86	0.12	17,17,17,17	0
57	MG	XA	1608	1/1	0.86	0.12	6,6,6,6	0
57	MG	YA	3189	1/1	0.86	0.14	27,27,27,27	0
57	MG	YA	3275	1/1	0.86	0.27	66,66,66,66	0
57	MG	RA	3181	1/1	0.86	0.19	28,28,28,28	0
57	MG	YA	3314	1/1	0.86	0.26	47,47,47,47	0
57	MG	RA	3273	1/1	0.86	0.32	45,45,45,45	0
57	MG	QA	1688	1/1	0.86	0.21	58,58,58,58	0
57	MG	XA	1665	1/1	0.86	0.08	72,72,72,72	0
57	MG	YA	3210	1/1	0.86	0.14	33,33,33,33	0
57	MG	RA	3008	1/1	0.86	0.15	16,16,16,16	0
57	MG	YA	3152	1/1	0.86	0.10	15,15,15,15	0
57	MG	YA	3154	1/1	0.86	0.30	56,56,56,56	0
57	MG	QA	1651	1/1	0.86	0.14	56,56,56,56	0
57	MG	YA	3334	1/1	0.86	0.21	48,48,48,48	0
57	MG	RA	3201	1/1	0.86	0.19	45,45,45,45	0
57	MG	RA	3296	1/1	0.86	0.16	40,40,40,40	0
57	MG	YA	3054	1/1	0.86	0.18	34,34,34,34	0
57	MG	XA	1701	1/1	0.86	0.13	26,26,26,26	0
57	MG	YP	202	1/1	0.86	0.40	29,29,29,29	0
57	MG	RA	3299	1/1	0.86	0.22	45,45,45,45	0
57	MG	YA	3197	1/1	0.87	0.31	43,43,43,43	0
57	MG	QA	1663	1/1	0.87	0.28	27,27,27,27	0
57	MG	YA	3132	1/1	0.87	0.22	35,35,35,35	0
57	MG	XA	1620	1/1	0.87	0.27	22,22,22,22	0
57	MG	QA	1689	1/1	0.87	0.12	78,78,78,78	0
57	MG	QA	1636	1/1	0.87	0.26	39,39,39,39	0
57	MG	RA	3283	1/1	0.87	0.26	42,42,42,42	0
57	MG	QA	1670	1/1	0.87	0.13	25,25,25,25	0
57	MG	QA	1673	1/1	0.87	0.15	35,35,35,35	0
57	MG	QA	1613	1/1	0.87	0.17	26,26,26,26	0
57	MG	RA	3177	1/1	0.87	0.20	25,25,25,25	0
57	MG	YA	3330	1/1	0.87	0.23	52,52,52,52	0
57	MG	YA	3333	1/1	0.87	0.30	43,43,43,43	0
57	MG	QA	1603	1/1	0.87	0.11	9,9,9,9	0
57	MG	XA	1706	1/1	0.87	0.19	47,47,47,47	0
57	MG	YA	3107	1/1	0.87	0.17	1,1,1,1	0
57	MG	RA	3272	1/1	0.87	0.26	48,48,48,48	0
57	MG	RA	3141	1/1	0.87	0.28	5,5,5,5	0
57	MG	RA	3147	1/1	0.87	0.16	28,28,28,28	0
57	MG	YA	3124	1/1	0.87	0.22	32,32,32,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	QA	1645	1/1	0.88	0.21	41,41,41,41	0
57	MG	XA	1622	1/1	0.88	0.18	33,33,33,33	0
57	MG	XA	1628	1/1	0.88	0.12	36,36,36,36	0
57	MG	XA	1651	1/1	0.88	0.28	50,50,50,50	0
57	MG	YA	3195	1/1	0.88	0.13	34,34,34,34	0
57	MG	XA	1652	1/1	0.88	0.16	44,44,44,44	0
57	MG	RA	3105	1/1	0.88	0.12	14,14,14,14	0
57	MG	RA	3298	1/1	0.88	0.20	56,56,56,56	0
57	MG	RA	3168	1/1	0.88	0.19	38,38,38,38	0
57	MG	RA	3300	1/1	0.88	0.29	43,43,43,43	0
57	MG	RA	3170	1/1	0.88	0.10	24,24,24,24	0
57	MG	RA	3112	1/1	0.88	0.18	31,31,31,31	0
57	MG	RA	3058	1/1	0.88	0.13	1,1,1,1	0
57	MG	RA	3213	1/1	0.88	0.15	38,38,38,38	0
57	MG	XA	1702	1/1	0.88	0.19	38,38,38,38	0
57	MG	YA	3223	1/1	0.88	0.19	27,27,27,27	0
57	MG	YA	3139	1/1	0.88	0.16	46,46,46,46	0
57	MG	YA	3143	1/1	0.88	0.15	25,25,25,25	0
57	MG	RA	3001	1/1	0.88	0.20	7,7,7,7	0
57	MG	RA	3005	1/1	0.88	0.25	4,4,4,4	0
57	MG	XA	1710	1/1	0.88	0.16	29,29,29,29	0
57	MG	RA	3184	1/1	0.88	0.18	35,35,35,35	0
57	MG	YA	3329	1/1	0.88	0.13	51,51,51,51	0
57	MG	RA	3007	1/1	0.88	0.28	22,22,22,22	0
57	MG	YA	3158	1/1	0.88	0.09	18,18,18,18	0
57	MG	YA	3161	1/1	0.88	0.19	32,32,32,32	0
57	MG	RA	3289	1/1	0.88	0.30	54,54,54,54	0
57	MG	R1	101	1/1	0.88	0.14	32,32,32,32	0
57	MG	RA	3151	1/1	0.88	0.11	13,13,13,13	0
57	MG	RA	3247	1/1	0.88	0.30	58,58,58,58	0
57	MG	YA	3178	1/1	0.88	0.28	40,40,40,40	0
57	MG	QA	1639	1/1	0.88	0.15	27,27,27,27	0
57	MG	YA	3188	1/1	0.89	0.17	31,31,31,31	0
57	MG	QA	1656	1/1	0.89	0.13	39,39,39,39	0
57	MG	XA	1683	1/1	0.89	0.11	20,20,20,20	0
57	MG	YA	3083	1/1	0.89	0.15	15,15,15,15	0
57	MG	RA	3219	1/1	0.89	0.12	31,31,31,31	0
57	MG	RA	3059	1/1	0.89	0.22	1,1,1,1	0
57	MG	XA	1688	1/1	0.89	0.21	30,30,30,30	0
57	MG	QA	1604	1/1	0.89	0.25	5,5,5,5	0
57	MG	RA	3241	1/1	0.89	0.10	51,51,51,51	0
57	MG	QA	1666	1/1	0.89	0.28	32,32,32,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3129	1/1	0.89	0.12	16,16,16,16	0
57	MG	QA	1617	1/1	0.89	0.20	15,15,15,15	0
57	MG	RA	3292	1/1	0.89	0.29	45,45,45,45	0
57	MG	XA	1708	1/1	0.89	0.17	43,43,43,43	0
57	MG	RA	3200	1/1	0.89	0.08	31,31,31,31	0
57	MG	RA	3294	1/1	0.89	0.21	54,54,54,54	0
57	MG	YA	3319	1/1	0.89	0.19	56,56,56,56	0
57	MG	QA	1622	1/1	0.89	0.14	42,42,42,42	0
57	MG	YA	3250	1/1	0.89	0.13	50,50,50,50	0
57	MG	QA	1612	1/1	0.89	0.31	22,22,22,22	0
57	MG	RA	3099	1/1	0.89	0.21	9,9,9,9	0
57	MG	XA	1645	1/1	0.89	0.20	25,25,25,25	0
57	MG	YA	3270	1/1	0.89	0.20	20,20,20,20	0
57	MG	RA	3025	1/1	0.89	0.14	3,3,3,3	0
57	MG	RA	3047	1/1	0.89	0.23	6,6,6,6	0
57	MG	YA	3024	1/1	0.89	0.30	3,3,3,3	0
57	MG	RA	3276	1/1	0.89	0.25	51,51,51,51	0
57	MG	QA	1694	1/1	0.89	0.13	50,50,50,50	0
57	MG	YA	3040	1/1	0.89	0.31	1,1,1,1	0
57	MG	YA	3041	1/1	0.89	0.30	18,18,18,18	0
57	MG	XA	1661	1/1	0.89	0.14	22,22,22,22	0
57	MG	RA	3140	1/1	0.89	0.14	20,20,20,20	0
57	MG	YA	3059	1/1	0.89	0.20	38,38,38,38	0
57	MG	Y0	103	1/1	0.89	0.24	50,50,50,50	0
57	MG	XA	1687	1/1	0.90	0.10	54,54,54,54	0
57	MG	RA	3027	1/1	0.90	0.29	2,2,2,2	0
57	MG	XA	1691	1/1	0.90	0.11	34,34,34,34	0
57	MG	YA	3294	1/1	0.90	0.17	48,48,48,48	0
57	MG	RA	3248	1/1	0.90	0.11	32,32,32,32	0
57	MG	RA	3043	1/1	0.90	0.18	5,5,5,5	0
57	MG	QA	1643	1/1	0.90	0.12	28,28,28,28	0
57	MG	XA	1705	1/1	0.90	0.23	58,58,58,58	0
57	MG	YA	3125	1/1	0.90	0.15	14,14,14,14	0
57	MG	RA	3004	1/1	0.90	0.19	11,11,11,11	0
57	MG	QA	1627	1/1	0.90	0.10	37,37,37,37	0
57	MG	QA	1648	1/1	0.90	0.10	49,49,49,49	0
57	MG	RA	3206	1/1	0.90	0.25	45,45,45,45	0
57	MG	QA	1693	1/1	0.90	0.20	49,49,49,49	0
57	MG	RA	3157	1/1	0.90	0.15	21,21,21,21	0
57	MG	QA	1676	1/1	0.90	0.16	48,48,48,48	0
57	MG	YA	3236	1/1	0.90	0.27	6,6,6,6	0
57	MG	QA	1677	1/1	0.90	0.19	26,26,26,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3248	1/1	0.90	0.28	37,37,37,37	0
57	MG	RA	3212	1/1	0.90	0.12	55,55,55,55	0
57	MG	RA	3165	1/1	0.90	0.15	19,19,19,19	0
57	MG	QE	201	1/1	0.90	0.17	37,37,37,37	0
57	MG	YA	3264	1/1	0.90	0.15	40,40,40,40	0
57	MG	RA	3086	1/1	0.90	0.21	5,5,5,5	0
57	MG	YA	3159	1/1	0.90	0.23	38,38,38,38	0
57	MG	RA	3088	1/1	0.90	0.24	12,12,12,12	0
57	MG	YA	3332	1/1	0.90	0.15	32,32,32,32	0
57	MG	RA	3225	1/1	0.90	0.19	28,28,28,28	0
57	MG	QA	1664	1/1	0.90	0.22	31,31,31,31	0
57	MG	RA	3019	1/1	0.90	0.16	8,8,8,8	0
57	MG	XA	1679	1/1	0.90	0.28	43,43,43,43	0
57	MG	YA	3053	1/1	0.90	0.22	3,3,3,3	0
57	MG	RA	3305	1/1	0.90	0.31	41,41,41,41	0
57	MG	QA	1665	1/1	0.90	0.09	18,18,18,18	0
57	MG	QA	1618	1/1	0.90	0.17	20,20,20,20	0
57	MG	YA	3072	1/1	0.90	0.14	2,2,2,2	0
57	MG	QA	1686	1/1	0.91	0.19	38,38,38,38	0
57	MG	YA	3110	1/1	0.91	0.08	11,11,11,11	0
57	MG	XA	1660	1/1	0.91	0.12	44,44,44,44	0
57	MG	YA	3194	1/1	0.91	0.06	27,27,27,27	0
57	MG	YA	3117	1/1	0.91	0.18	12,12,12,12	0
57	MG	QA	1687	1/1	0.91	0.26	40,40,40,40	0
57	MG	QA	1674	1/1	0.91	0.18	26,26,26,26	0
57	MG	RA	3259	1/1	0.91	0.17	37,37,37,37	0
57	MG	RA	3260	1/1	0.91	0.19	49,49,49,49	0
57	MG	RA	3239	1/1	0.91	0.17	37,37,37,37	0
57	MG	YA	3006	1/1	0.91	0.31	1,1,1,1	0
57	MG	YA	3011	1/1	0.91	0.22	2,2,2,2	0
57	MG	YA	3014	1/1	0.91	0.23	3,3,3,3	0
57	MG	YA	3018	1/1	0.91	0.25	4,4,4,4	0
57	MG	YA	3142	1/1	0.91	0.23	39,39,39,39	0
57	MG	QA	1659	1/1	0.91	0.15	33,33,33,33	0
57	MG	YA	3224	1/1	0.91	0.14	13,13,13,13	0
57	MG	YA	3144	1/1	0.91	0.25	44,44,44,44	0
57	MG	YA	3318	1/1	0.91	0.21	38,38,38,38	0
57	MG	QA	1626	1/1	0.91	0.07	29,29,29,29	0
57	MG	QA	1624	1/1	0.91	0.10	39,39,39,39	0
57	MG	RA	3185	1/1	0.91	0.24	36,36,36,36	0
57	MG	YA	3153	1/1	0.91	0.20	31,31,31,31	0
57	MG	RA	3302	1/1	0.91	0.12	51,51,51,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	QA	1632	1/1	0.91	0.21	16,16,16,16	0
57	MG	YA	3052	1/1	0.91	0.15	14,14,14,14	0
57	MG	XA	1629	1/1	0.91	0.13	20,20,20,20	0
57	MG	XA	1632	1/1	0.91	0.20	18,18,18,18	0
57	MG	RA	3050	1/1	0.91	0.22	7,7,7,7	0
57	MG	RA	3167	1/1	0.91	0.10	38,38,38,38	0
57	MG	YA	3168	1/1	0.91	0.15	32,32,32,32	0
57	MG	RA	3144	1/1	0.91	0.12	32,32,32,32	0
57	MG	XA	1653	1/1	0.91	0.20	26,26,26,26	0
57	MG	YA	3077	1/1	0.91	0.19	11,11,11,11	0
57	MG	YA	3281	1/1	0.91	0.18	44,44,44,44	0
57	MG	YD	301	1/1	0.91	0.20	16,16,16,16	0
57	MG	YE	301	1/1	0.91	0.17	2,2,2,2	0
57	MG	YA	3080	1/1	0.91	0.26	9,9,9,9	0
57	MG	RB	201	1/1	0.91	0.10	59,59,59,59	0
57	MG	YA	3094	1/1	0.91	0.17	9,9,9,9	0
58	PAR	QA	1696	42/42	0.91	0.15	61,61,61,62	0
57	MG	YA	3226	1/1	0.92	0.15	44,44,44,44	0
57	MG	YA	3229	1/1	0.92	0.14	50,50,50,50	0
57	MG	QA	1653	1/1	0.92	0.20	61,61,61,61	0
57	MG	YA	3234	1/1	0.92	0.28	54,54,54,54	0
57	MG	YA	3084	1/1	0.92	0.14	18,18,18,18	0
57	MG	RA	3074	1/1	0.92	0.18	2,2,2,2	0
57	MG	YA	3098	1/1	0.92	0.21	12,12,12,12	0
57	MG	YA	3102	1/1	0.92	0.12	9,9,9,9	0
57	MG	RA	3156	1/1	0.92	0.11	29,29,29,29	0
57	MG	RA	3076	1/1	0.92	0.16	14,14,14,14	0
57	MG	YA	3111	1/1	0.92	0.18	25,25,25,25	0
57	MG	RA	3029	1/1	0.92	0.14	3,3,3,3	0
57	MG	RA	3032	1/1	0.92	0.25	28,28,28,28	0
57	MG	RA	3162	1/1	0.92	0.06	46,46,46,46	0
57	MG	XA	1689	1/1	0.92	0.16	21,21,21,21	0
57	MG	RA	3164	1/1	0.92	0.05	28,28,28,28	0
57	MG	QA	1691	1/1	0.92	0.18	26,26,26,26	0
57	MG	XA	1696	1/1	0.92	0.15	25,25,25,25	0
57	MG	RA	3044	1/1	0.92	0.15	22,22,22,22	0
57	MG	YA	3131	1/1	0.92	0.10	35,35,35,35	0
57	MG	YA	3283	1/1	0.92	0.15	34,34,34,34	0
57	MG	QA	1683	1/1	0.92	0.16	37,37,37,37	0
57	MG	QA	1631	1/1	0.92	0.07	32,32,32,32	0
57	MG	RA	3232	1/1	0.92	0.17	1,1,1,1	0
57	MG	RA	3002	1/1	0.92	0.19	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	R5	101	1/1	0.92	0.08	12,12,12,12	0
57	MG	XA	1709	1/1	0.92	0.30	49,49,49,49	0
57	MG	RA	3118	1/1	0.92	0.15	35,35,35,35	0
57	MG	XA	1602	1/1	0.92	0.27	8,8,8,8	0
57	MG	RA	3124	1/1	0.92	0.20	12,12,12,12	0
57	MG	XA	1614	1/1	0.92	0.19	4,4,4,4	0
57	MG	RA	3242	1/1	0.92	0.31	46,46,46,46	0
57	MG	RA	3179	1/1	0.92	0.07	15,15,15,15	0
57	MG	XA	1621	1/1	0.92	0.09	30,30,30,30	0
57	MG	RA	3180	1/1	0.92	0.24	40,40,40,40	0
57	MG	YA	3004	1/1	0.92	0.23	7,7,7,7	0
57	MG	XA	1627	1/1	0.92	0.08	4,4,4,4	0
57	MG	RA	3131	1/1	0.92	0.24	11,11,11,11	0
57	MG	YA	3310	1/1	0.92	0.12	32,32,32,32	0
57	MG	RA	3132	1/1	0.92	0.11	46,46,46,46	0
57	MG	YA	3312	1/1	0.92	0.14	33,33,33,33	0
57	MG	RA	3057	1/1	0.92	0.15	9,9,9,9	0
57	MG	YA	3023	1/1	0.92	0.23	15,15,15,15	0
57	MG	XA	1633	1/1	0.92	0.12	28,28,28,28	0
57	MG	XA	1634	1/1	0.92	0.34	27,27,27,27	0
57	MG	XA	1635	1/1	0.92	0.20	11,11,11,11	0
57	MG	YA	3032	1/1	0.92	0.26	1,1,1,1	0
57	MG	YA	3034	1/1	0.92	0.26	10,10,10,10	0
57	MG	YA	3038	1/1	0.92	0.17	24,24,24,24	0
57	MG	XA	1636	1/1	0.92	0.13	52,52,52,52	0
57	MG	QA	1647	1/1	0.92	0.20	65,65,65,65	0
57	MG	YA	3044	1/1	0.92	0.27	2,2,2,2	0
57	MG	XA	1646	1/1	0.92	0.19	36,36,36,36	0
57	MG	YA	3200	1/1	0.92	0.11	33,33,33,33	0
57	MG	RA	3189	1/1	0.92	0.08	12,12,12,12	0
57	MG	YA	3205	1/1	0.92	0.09	56,56,56,56	0
57	MG	RA	3022	1/1	0.92	0.19	15,15,15,15	0
57	MG	RA	3254	1/1	0.92	0.21	60,60,60,60	0
57	MG	YA	3055	1/1	0.92	0.16	9,9,9,9	0
57	MG	XA	1655	1/1	0.92	0.17	21,21,21,21	0
57	MG	YA	3215	1/1	0.92	0.11	25,25,25,25	0
57	MG	RA	3197	1/1	0.92	0.17	26,26,26,26	0
57	MG	YA	3071	1/1	0.92	0.09	14,14,14,14	0
57	MG	YP	201	1/1	0.92	0.12	22,22,22,22	0
57	MG	RA	3256	1/1	0.92	0.26	33,33,33,33	0
57	MG	QA	1629	1/1	0.92	0.19	69,69,69,69	0
57	MG	Y0	102	1/1	0.92	0.22	27,27,27,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3066	1/1	0.92	0.23	1,1,1,1	0
57	MG	QA	1614	1/1	0.92	0.18	29,29,29,29	0
57	MG	RA	3060	1/1	0.93	0.19	1,1,1,1	0
57	MG	YA	3135	1/1	0.93	0.23	30,30,30,30	0
57	MG	XA	1654	1/1	0.93	0.11	42,42,42,42	0
57	MG	RA	3172	1/1	0.93	0.12	26,26,26,26	0
57	MG	YA	3268	1/1	0.93	0.13	17,17,17,17	0
57	MG	RA	3223	1/1	0.93	0.12	7,7,7,7	0
57	MG	RA	3174	1/1	0.93	0.11	20,20,20,20	0
57	MG	YA	3273	1/1	0.93	0.20	37,37,37,37	0
57	MG	RA	3127	1/1	0.93	0.18	5,5,5,5	0
57	MG	RA	3128	1/1	0.93	0.11	12,12,12,12	0
57	MG	YA	3149	1/1	0.93	0.21	22,22,22,22	0
57	MG	RA	3129	1/1	0.93	0.18	38,38,38,38	0
57	MG	YA	3279	1/1	0.93	0.19	30,30,30,30	0
57	MG	RA	3028	1/1	0.93	0.13	2,2,2,2	0
57	MG	YA	3039	1/1	0.93	0.13	2,2,2,2	0
57	MG	QA	1685	1/1	0.93	0.14	38,38,38,38	0
57	MG	QA	1672	1/1	0.93	0.09	24,24,24,24	0
57	MG	YA	3043	1/1	0.93	0.16	2,2,2,2	0
57	MG	YA	3286	1/1	0.93	0.13	46,46,46,46	0
57	MG	RA	3280	1/1	0.93	0.20	47,47,47,47	0
57	MG	RA	3039	1/1	0.93	0.15	1,1,1,1	0
57	MG	QA	1623	1/1	0.93	0.19	26,26,26,26	0
57	MG	RA	3284	1/1	0.93	0.31	32,32,32,32	0
57	MG	XA	1613	1/1	0.93	0.13	18,18,18,18	0
57	MG	RA	3015	1/1	0.93	0.08	1,1,1,1	0
57	MG	QA	1681	1/1	0.93	0.19	43,43,43,43	0
57	MG	RA	3195	1/1	0.93	0.11	4,4,4,4	0
57	MG	YA	3069	1/1	0.93	0.15	15,15,15,15	0
57	MG	YA	3184	1/1	0.93	0.09	32,32,32,32	0
57	MG	XA	1699	1/1	0.93	0.16	44,44,44,44	0
57	MG	RA	3021	1/1	0.93	0.21	1,1,1,1	0
57	MG	QA	1657	1/1	0.93	0.07	55,55,55,55	0
57	MG	XA	1626	1/1	0.93	0.13	23,23,23,23	0
57	MG	YA	3190	1/1	0.93	0.17	6,6,6,6	0
57	MG	YA	3079	1/1	0.93	0.34	10,10,10,10	0
57	MG	RA	3054	1/1	0.93	0.21	5,5,5,5	0
57	MG	RA	3098	1/1	0.93	0.26	1,1,1,1	0
57	MG	YA	3313	1/1	0.93	0.17	23,23,23,23	0
57	MG	YA	3196	1/1	0.93	0.26	34,34,34,34	0
57	MG	RA	3055	1/1	0.93	0.15	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3085	1/1	0.93	0.23	6,6,6,6	0
57	MG	YA	3088	1/1	0.93	0.20	5,5,5,5	0
57	MG	YA	3089	1/1	0.93	0.15	25,25,25,25	0
57	MG	XA	1630	1/1	0.93	0.13	27,27,27,27	0
57	MG	XA	1631	1/1	0.93	0.17	12,12,12,12	0
57	MG	QF	201	1/1	0.93	0.08	41,41,41,41	0
57	MG	YA	3104	1/1	0.93	0.14	26,26,26,26	0
57	MG	YA	3106	1/1	0.93	0.17	1,1,1,1	0
57	MG	YA	3327	1/1	0.93	0.19	56,56,56,56	0
57	MG	RA	3109	1/1	0.93	0.16	23,23,23,23	0
57	MG	RA	3111	1/1	0.93	0.18	22,22,22,22	0
57	MG	QA	1615	1/1	0.93	0.10	65,65,65,65	0
57	MG	YA	3331	1/1	0.93	0.22	43,43,43,43	0
57	MG	RA	3117	1/1	0.93	0.08	43,43,43,43	0
57	MG	XA	1644	1/1	0.93	0.12	27,27,27,27	0
57	MG	QA	1684	1/1	0.93	0.11	40,40,40,40	0
57	MG	YA	3225	1/1	0.93	0.12	23,23,23,23	0
57	MG	YA	3002	1/1	0.93	0.22	1,1,1,1	0
57	MG	YA	3003	1/1	0.93	0.18	3,3,3,3	0
57	MG	RA	3119	1/1	0.93	0.26	43,43,43,43	0
57	MG	YB	203	1/1	0.93	0.20	20,20,20,20	0
57	MG	YA	3127	1/1	0.93	0.20	12,12,12,12	0
57	MG	YA	3235	1/1	0.93	0.15	18,18,18,18	0
57	MG	RA	3121	1/1	0.93	0.16	20,20,20,20	0
57	MG	YA	3237	1/1	0.93	0.27	3,3,3,3	0
57	MG	RA	3215	1/1	0.93	0.22	34,34,34,34	0
57	MG	YA	3246	1/1	0.93	0.40	39,39,39,39	0
57	MG	YA	3130	1/1	0.93	0.06	25,25,25,25	0
57	MG	YA	3013	1/1	0.93	0.19	4,4,4,4	0
58	PAR	XA	1715	42/42	0.93	0.12	52,52,53,53	0
57	MG	RA	3234	1/1	0.94	0.07	2,2,2,2	0
57	MG	YA	3138	1/1	0.94	0.22	9,9,9,9	0
57	MG	XA	1680	1/1	0.94	0.15	29,29,29,29	0
57	MG	YA	3252	1/1	0.94	0.21	23,23,23,23	0
57	MG	YA	3256	1/1	0.94	0.19	13,13,13,13	0
57	MG	RA	3101	1/1	0.94	0.13	14,14,14,14	0
57	MG	YA	3258	1/1	0.94	0.12	12,12,12,12	0
57	MG	YA	3262	1/1	0.94	0.20	5,5,5,5	0
57	MG	XA	1684	1/1	0.94	0.11	26,26,26,26	0
57	MG	RA	3104	1/1	0.94	0.09	4,4,4,4	0
57	MG	YA	3266	1/1	0.94	0.14	40,40,40,40	0
57	MG	RA	3279	1/1	0.94	0.25	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	QA	1658	1/1	0.94	0.08	78,78,78,78	0
57	MG	QA	1641	1/1	0.94	0.09	16,16,16,16	0
57	MG	YA	3151	1/1	0.94	0.12	10,10,10,10	0
57	MG	XA	1615	1/1	0.94	0.24	2,2,2,2	0
57	MG	XA	1690	1/1	0.94	0.10	48,48,48,48	0
57	MG	RA	3243	1/1	0.94	0.14	37,37,37,37	0
57	MG	YA	3276	1/1	0.94	0.11	7,7,7,7	0
57	MG	YA	3155	1/1	0.94	0.12	18,18,18,18	0
57	MG	YA	3058	1/1	0.94	0.18	7,7,7,7	0
57	MG	RA	3190	1/1	0.94	0.23	53,53,53,53	0
57	MG	YA	3062	1/1	0.94	0.18	5,5,5,5	0
57	MG	YA	3160	1/1	0.94	0.07	6,6,6,6	0
57	MG	QV	103	1/1	0.94	0.20	30,30,30,30	0
57	MG	YA	3067	1/1	0.94	0.14	31,31,31,31	0
57	MG	YA	3068	1/1	0.94	0.18	4,4,4,4	0
57	MG	RA	3153	1/1	0.94	0.17	19,19,19,19	0
57	MG	YA	3287	1/1	0.94	0.27	47,47,47,47	0
57	MG	YA	3170	1/1	0.94	0.05	29,29,29,29	0
57	MG	YA	3171	1/1	0.94	0.17	13,13,13,13	0
57	MG	YA	3070	1/1	0.94	0.09	5,5,5,5	0
57	MG	RA	3196	1/1	0.94	0.12	55,55,55,55	0
57	MG	YA	3175	1/1	0.94	0.25	34,34,34,34	0
57	MG	YA	3177	1/1	0.94	0.19	20,20,20,20	0
57	MG	RA	3069	1/1	0.94	0.08	27,27,27,27	0
57	MG	YA	3179	1/1	0.94	0.11	19,19,19,19	0
57	MG	XA	1703	1/1	0.94	0.09	25,25,25,25	0
57	MG	YA	3299	1/1	0.94	0.28	48,48,48,48	0
57	MG	YA	3181	1/1	0.94	0.13	15,15,15,15	0
57	MG	YA	3182	1/1	0.94	0.10	30,30,30,30	0
57	MG	YA	3076	1/1	0.94	0.15	14,14,14,14	0
57	MG	YA	3303	1/1	0.94	0.20	35,35,35,35	0
57	MG	RA	3048	1/1	0.94	0.10	1,1,1,1	0
57	MG	QA	1668	1/1	0.94	0.26	25,25,25,25	0
57	MG	QA	1678	1/1	0.94	0.23	41,41,41,41	0
57	MG	RA	3120	1/1	0.94	0.20	18,18,18,18	0
57	MG	RA	3081	1/1	0.94	0.11	11,11,11,11	0
57	MG	RA	3122	1/1	0.94	0.14	25,25,25,25	0
57	MG	YA	3193	1/1	0.94	0.21	18,18,18,18	0
57	MG	QA	1655	1/1	0.94	0.16	24,24,24,24	0
57	MG	YA	3315	1/1	0.94	0.21	17,17,17,17	0
57	MG	RA	3258	1/1	0.94	0.14	26,26,26,26	0
57	MG	YA	3093	1/1	0.94	0.16	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3209	1/1	0.94	0.19	33,33,33,33	0
57	MG	YA	3095	1/1	0.94	0.12	2,2,2,2	0
57	MG	QA	1652	1/1	0.94	0.08	91,91,91,91	0
57	MG	YA	3100	1/1	0.94	0.21	6,6,6,6	0
57	MG	RA	3087	1/1	0.94	0.17	35,35,35,35	0
57	MG	YA	3208	1/1	0.94	0.18	31,31,31,31	0
57	MG	YA	3324	1/1	0.94	0.13	28,28,28,28	0
57	MG	RA	3030	1/1	0.94	0.19	4,4,4,4	0
57	MG	XV	103	1/1	0.94	0.10	7,7,7,7	0
57	MG	RA	3003	1/1	0.94	0.16	3,3,3,3	0
57	MG	RA	3173	1/1	0.94	0.10	38,38,38,38	0
57	MG	RA	3092	1/1	0.94	0.17	12,12,12,12	0
57	MG	YA	3217	1/1	0.94	0.06	43,43,43,43	0
57	MG	RA	3134	1/1	0.94	0.06	22,22,22,22	0
57	MG	RA	3093	1/1	0.94	0.20	11,11,11,11	0
57	MG	RA	3139	1/1	0.94	0.18	17,17,17,17	0
57	MG	XA	1657	1/1	0.94	0.09	23,23,23,23	0
57	MG	RA	3224	1/1	0.94	0.12	11,11,11,11	0
57	MG	RA	3034	1/1	0.94	0.22	4,4,4,4	0
57	MG	RA	3228	1/1	0.94	0.17	8,8,8,8	0
57	MG	YA	3228	1/1	0.94	0.07	21,21,21,21	0
57	MG	QA	1642	1/1	0.94	0.06	37,37,37,37	0
57	MG	YA	3232	1/1	0.94	0.13	6,6,6,6	0
57	MG	YE	302	1/1	0.94	0.14	4,4,4,4	0
57	MG	RA	3233	1/1	0.94	0.11	18,18,18,18	0
57	MG	XA	1667	1/1	0.94	0.07	25,25,25,25	0
57	MG	XA	1674	1/1	0.94	0.22	16,16,16,16	0
57	MG	YA	3033	1/1	0.94	0.17	4,4,4,4	0
57	MG	YA	3134	1/1	0.94	0.14	17,17,17,17	0
57	MG	XA	1677	1/1	0.94	0.22	21,21,21,21	0
57	MG	YA	3243	1/1	0.94	0.22	18,18,18,18	0
59	ZN	XD	301	1/1	0.94	0.18	43,43,43,43	0
57	MG	RA	3090	1/1	0.95	0.13	9,9,9,9	0
57	MG	XA	1649	1/1	0.95	0.06	27,27,27,27	0
57	MG	YA	3259	1/1	0.95	0.08	14,14,14,14	0
57	MG	RA	3053	1/1	0.95	0.21	1,1,1,1	0
57	MG	RA	3186	1/1	0.95	0.12	34,34,34,34	0
57	MG	QA	1662	1/1	0.95	0.13	27,27,27,27	0
57	MG	RA	3245	1/1	0.95	0.28	57,57,57,57	0
57	MG	QA	1611	1/1	0.95	0.13	2,2,2,2	0
57	MG	RA	3297	1/1	0.95	0.08	29,29,29,29	0
57	MG	RA	3094	1/1	0.95	0.12	13,13,13,13	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3191	1/1	0.95	0.08	36,36,36,36	0
57	MG	RA	3095	1/1	0.95	0.22	7,7,7,7	0
57	MG	QA	1601	1/1	0.95	0.17	38,38,38,38	0
57	MG	YA	3046	1/1	0.95	0.20	3,3,3,3	0
57	MG	RA	3150	1/1	0.95	0.10	15,15,15,15	0
57	MG	QA	1644	1/1	0.95	0.10	33,33,33,33	0
57	MG	QA	1602	1/1	0.95	0.26	18,18,18,18	0
57	MG	RA	3033	1/1	0.95	0.27	4,4,4,4	0
57	MG	RA	3154	1/1	0.95	0.19	23,23,23,23	0
57	MG	XA	1678	1/1	0.95	0.18	23,23,23,23	0
57	MG	YA	3172	1/1	0.95	0.05	44,44,44,44	0
57	MG	YA	3060	1/1	0.95	0.19	5,5,5,5	0
57	MG	RA	3155	1/1	0.95	0.12	11,11,11,11	0
57	MG	YA	3064	1/1	0.95	0.19	12,12,12,12	0
57	MG	YA	3176	1/1	0.95	0.25	41,41,41,41	0
57	MG	RA	3203	1/1	0.95	0.04	27,27,27,27	0
57	MG	XA	1681	1/1	0.95	0.22	23,23,23,23	0
57	MG	RA	3061	1/1	0.95	0.17	1,1,1,1	0
57	MG	RB	204	1/1	0.95	0.30	39,39,39,39	0
57	MG	QA	1680	1/1	0.95	0.07	18,18,18,18	0
57	MG	RA	3016	1/1	0.95	0.20	11,11,11,11	0
57	MG	RA	3160	1/1	0.95	0.16	43,43,43,43	0
57	MG	RA	3040	1/1	0.95	0.18	1,1,1,1	0
57	MG	RA	3042	1/1	0.95	0.13	23,23,23,23	0
57	MG	YA	3298	1/1	0.95	0.07	16,16,16,16	0
57	MG	RA	3163	1/1	0.95	0.21	43,43,43,43	0
57	MG	RA	3070	1/1	0.95	0.17	4,4,4,4	0
57	MG	XA	1605	1/1	0.95	0.29	26,26,26,26	0
57	MG	YA	3191	1/1	0.95	0.07	34,34,34,34	0
57	MG	YA	3081	1/1	0.95	0.13	14,14,14,14	0
57	MG	QA	1637	1/1	0.95	0.15	43,43,43,43	0
57	MG	XA	1697	1/1	0.95	0.07	19,19,19,19	0
57	MG	XA	1611	1/1	0.95	0.14	11,11,11,11	0
57	MG	YA	3309	1/1	0.95	0.33	57,57,57,57	0
57	MG	RA	3020	1/1	0.95	0.20	1,1,1,1	0
57	MG	RA	3216	1/1	0.95	0.17	4,4,4,4	0
57	MG	RA	3045	1/1	0.95	0.17	11,11,11,11	0
57	MG	YA	3202	1/1	0.95	0.05	15,15,15,15	0
57	MG	XA	1617	1/1	0.95	0.20	4,4,4,4	0
57	MG	RA	3271	1/1	0.95	0.25	41,41,41,41	0
57	MG	XA	1619	1/1	0.95	0.11	46,46,46,46	0
57	MG	YA	3207	1/1	0.95	0.21	34,34,34,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3080	1/1	0.95	0.19	20,20,20,20	0
57	MG	RA	3222	1/1	0.95	0.18	18,18,18,18	0
57	MG	YA	3103	1/1	0.95	0.07	15,15,15,15	0
57	MG	RA	3123	1/1	0.95	0.14	34,34,34,34	0
57	MG	XA	1623	1/1	0.95	0.17	14,14,14,14	0
57	MG	RA	3046	1/1	0.95	0.13	11,11,11,11	0
57	MG	YA	3109	1/1	0.95	0.16	23,23,23,23	0
57	MG	RA	3082	1/1	0.95	0.12	30,30,30,30	0
57	MG	RA	3226	1/1	0.95	0.15	32,32,32,32	0
57	MG	YA	3114	1/1	0.95	0.10	7,7,7,7	0
57	MG	YA	3222	1/1	0.95	0.07	25,25,25,25	0
57	MG	QA	1608	1/1	0.95	0.07	45,45,45,45	0
57	MG	RA	3229	1/1	0.95	0.22	3,3,3,3	0
57	MG	XV	102	1/1	0.95	0.17	10,10,10,10	0
57	MG	RA	3231	1/1	0.95	0.11	14,14,14,14	0
57	MG	YA	3227	1/1	0.95	0.12	19,19,19,19	0
57	MG	QA	1650	1/1	0.95	0.12	16,16,16,16	0
57	MG	QA	1610	1/1	0.95	0.20	13,13,13,13	0
57	MG	QA	1661	1/1	0.95	0.11	5,5,5,5	0
57	MG	RA	3236	1/1	0.95	0.09	13,13,13,13	0
57	MG	RA	3089	1/1	0.95	0.15	17,17,17,17	0
57	MG	XA	1638	1/1	0.95	0.10	46,46,46,46	0
57	MG	YA	3012	1/1	0.95	0.17	4,4,4,4	0
57	MG	XA	1639	1/1	0.95	0.19	6,6,6,6	0
57	MG	XA	1642	1/1	0.95	0.10	9,9,9,9	0
57	MG	YA	3239	1/1	0.95	0.13	10,10,10,10	0
57	MG	YQ	201	1/1	0.95	0.15	152,152,152,152	0
57	MG	YA	3016	1/1	0.95	0.09	2,2,2,2	0
57	MG	XA	1643	1/1	0.95	0.19	43,43,43,43	0
57	MG	RA	3238	1/1	0.95	0.21	43,43,43,43	0
57	MG	Y7	101	1/1	0.95	0.26	31,31,31,31	0
57	MG	RA	3137	1/1	0.95	0.09	3,3,3,3	0
57	MG	YA	3025	1/1	0.95	0.19	6,6,6,6	0
57	MG	YA	3026	1/1	0.95	0.25	5,5,5,5	0
57	MG	RA	3056	1/1	0.96	0.16	3,3,3,3	0
57	MG	YA	3271	1/1	0.96	0.24	35,35,35,35	0
57	MG	XA	1606	1/1	0.96	0.12	17,17,17,17	0
57	MG	RA	3176	1/1	0.96	0.07	26,26,26,26	0
57	MG	YA	3035	1/1	0.96	0.14	2,2,2,2	0
57	MG	YA	3105	1/1	0.96	0.09	5,5,5,5	0
57	MG	YA	3183	1/1	0.96	0.17	21,21,21,21	0
57	MG	YA	3037	1/1	0.96	0.20	3,3,3,3	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	XA	1647	1/1	0.96	0.10	53,53,53,53	0
57	MG	QV	102	1/1	0.96	0.15	11,11,11,11	0
57	MG	XA	1650	1/1	0.96	0.15	34,34,34,34	0
57	MG	RA	3110	1/1	0.96	0.18	11,11,11,11	0
57	MG	YA	3112	1/1	0.96	0.12	1,1,1,1	0
57	MG	YA	3113	1/1	0.96	0.19	4,4,4,4	0
57	MG	YA	3042	1/1	0.96	0.19	12,12,12,12	0
57	MG	QA	1675	1/1	0.96	0.07	32,32,32,32	0
57	MG	RA	3148	1/1	0.96	0.08	22,22,22,22	0
57	MG	RA	3227	1/1	0.96	0.11	14,14,14,14	0
57	MG	RA	3204	1/1	0.96	0.09	16,16,16,16	0
57	MG	YA	3121	1/1	0.96	0.14	19,19,19,19	0
57	MG	YA	3198	1/1	0.96	0.26	86,86,86,86	0
57	MG	YA	3123	1/1	0.96	0.09	4,4,4,4	0
57	MG	YA	3048	1/1	0.96	0.15	2,2,2,2	0
57	MG	YA	3049	1/1	0.96	0.26	2,2,2,2	0
57	MG	YA	3204	1/1	0.96	0.11	25,25,25,25	0
57	MG	YA	3050	1/1	0.96	0.21	2,2,2,2	0
57	MG	YA	3051	1/1	0.96	0.19	3,3,3,3	0
57	MG	RA	3183	1/1	0.96	0.06	37,37,37,37	0
57	MG	RA	3230	1/1	0.96	0.15	4,4,4,4	0
57	MG	RA	3304	1/1	0.96	0.22	29,29,29,29	0
57	MG	RA	3278	1/1	0.96	0.20	38,38,38,38	0
57	MG	QA	1671	1/1	0.96	0.07	25,25,25,25	0
57	MG	YA	3304	1/1	0.96	0.21	31,31,31,31	0
57	MG	YA	3214	1/1	0.96	0.14	4,4,4,4	0
57	MG	YA	3306	1/1	0.96	0.38	66,66,66,66	0
57	MG	XA	1624	1/1	0.96	0.06	18,18,18,18	0
57	MG	XA	1625	1/1	0.96	0.14	16,16,16,16	0
57	MG	YA	3061	1/1	0.96	0.18	4,4,4,4	0
57	MG	RA	3130	1/1	0.96	0.18	30,30,30,30	0
57	MG	YA	3219	1/1	0.96	0.27	37,37,37,37	0
57	MG	YA	3141	1/1	0.96	0.08	29,29,29,29	0
57	MG	RA	3166	1/1	0.96	0.07	9,9,9,9	0
57	MG	XA	1675	1/1	0.96	0.23	35,35,35,35	0
57	MG	XX	102	1/1	0.96	0.19	14,14,14,14	0
57	MG	YA	3147	1/1	0.96	0.11	19,19,19,19	0
57	MG	QX	102	1/1	0.96	0.24	45,45,45,45	0
57	MG	RA	3235	1/1	0.96	0.13	8,8,8,8	0
57	MG	RA	3096	1/1	0.96	0.12	11,11,11,11	0
57	MG	RA	3073	1/1	0.96	0.09	17,17,17,17	0
57	MG	RA	3136	1/1	0.96	0.20	2,2,2,2	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	YA	3231	1/1	0.96	0.18	4,4,4,4	0
57	MG	RE	302	1/1	0.96	0.10	8,8,8,8	0
57	MG	RF	301	1/1	0.96	0.08	52,52,52,52	0
57	MG	YA	3325	1/1	0.96	0.09	47,47,47,47	0
57	MG	RR	201	1/1	0.96	0.17	9,9,9,9	0
57	MG	YA	3078	1/1	0.96	0.15	12,12,12,12	0
57	MG	YA	3157	1/1	0.96	0.05	9,9,9,9	0
57	MG	YA	3015	1/1	0.96	0.22	11,11,11,11	0
57	MG	RA	3193	1/1	0.96	0.10	49,49,49,49	0
57	MG	YA	3017	1/1	0.96	0.15	13,13,13,13	0
57	MG	YA	3240	1/1	0.96	0.21	5,5,5,5	0
57	MG	YA	3242	1/1	0.96	0.27	1,1,1,1	0
57	MG	XA	1637	1/1	0.96	0.08	44,44,44,44	0
57	MG	YA	3020	1/1	0.96	0.18	5,5,5,5	0
57	MG	YA	3336	1/1	0.96	0.12	29,29,29,29	0
57	MG	YA	3163	1/1	0.96	0.11	34,34,34,34	0
57	MG	YA	3249	1/1	0.96	0.11	3,3,3,3	0
57	MG	YA	3166	1/1	0.96	0.26	36,36,36,36	0
57	MG	YA	3022	1/1	0.96	0.16	10,10,10,10	0
57	MG	YA	3087	1/1	0.96	0.16	4,4,4,4	0
57	MG	YA	3169	1/1	0.96	0.12	18,18,18,18	0
57	MG	RA	3240	1/1	0.96	0.20	20,20,20,20	0
57	MG	RA	3031	1/1	0.96	0.10	5,5,5,5	0
57	MG	YA	3260	1/1	0.96	0.19	15,15,15,15	0
57	MG	YA	3261	1/1	0.96	0.15	11,11,11,11	0
57	MG	YA	3092	1/1	0.96	0.23	9,9,9,9	0
57	MG	XA	1641	1/1	0.96	0.07	7,7,7,7	0
57	MG	RA	3075	1/1	0.96	0.06	41,41,41,41	0
57	MG	Y5	101	1/1	0.96	0.15	18,18,18,18	0
57	MG	XA	1693	1/1	0.96	0.23	36,36,36,36	0
57	MG	YA	3097	1/1	0.96	0.22	8,8,8,8	0
57	MG	RA	3041	1/1	0.96	0.18	4,4,4,4	0
57	MG	YA	3269	1/1	0.96	0.18	22,22,22,22	0
57	MG	RA	3125	1/1	0.97	0.17	20,20,20,20	0
57	MG	RA	3282	1/1	0.97	0.27	51,51,51,51	0
57	MG	YA	3036	1/1	0.97	0.10	2,2,2,2	0
57	MG	RA	3126	1/1	0.97	0.10	9,9,9,9	0
57	MG	RA	3084	1/1	0.97	0.09	13,13,13,13	0
57	MG	YA	3099	1/1	0.97	0.16	3,3,3,3	0
57	MG	RA	3205	1/1	0.97	0.20	19,19,19,19	0
57	MG	YA	3164	1/1	0.97	0.08	24,24,24,24	0
57	MG	YA	3165	1/1	0.97	0.15	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3085	1/1	0.97	0.13	4,4,4,4	0
57	MG	RE	301	1/1	0.97	0.15	2,2,2,2	0
57	MG	XA	1663	1/1	0.97	0.21	37,37,37,37	0
57	MG	XA	1664	1/1	0.97	0.14	18,18,18,18	0
57	MG	QA	1660	1/1	0.97	0.14	10,10,10,10	0
57	MG	QA	1633	1/1	0.97	0.08	25,25,25,25	0
57	MG	RA	3071	1/1	0.97	0.17	19,19,19,19	0
57	MG	XA	1668	1/1	0.97	0.08	32,32,32,32	0
57	MG	XA	1669	1/1	0.97	0.11	11,11,11,11	0
57	MG	XA	1670	1/1	0.97	0.10	14,14,14,14	0
57	MG	XA	1671	1/1	0.97	0.17	5,5,5,5	0
57	MG	QA	1646	1/1	0.97	0.08	56,56,56,56	0
57	MG	QA	1669	1/1	0.97	0.24	38,38,38,38	0
57	MG	XA	1676	1/1	0.97	0.13	15,15,15,15	0
57	MG	YA	3247	1/1	0.97	0.18	3,3,3,3	0
57	MG	YA	3001	1/1	0.97	0.14	3,3,3,3	0
57	MG	YA	3056	1/1	0.97	0.13	5,5,5,5	0
57	MG	RA	3159	1/1	0.97	0.06	28,28,28,28	0
57	MG	YA	3122	1/1	0.97	0.18	5,5,5,5	0
57	MG	RA	3135	1/1	0.97	0.07	20,20,20,20	0
57	MG	RA	3114	1/1	0.97	0.17	17,17,17,17	0
57	MG	XA	1604	1/1	0.97	0.20	19,19,19,19	0
57	MG	YA	3126	1/1	0.97	0.18	5,5,5,5	0
57	MG	YA	3007	1/1	0.97	0.12	4,4,4,4	0
57	MG	YA	3009	1/1	0.97	0.19	3,3,3,3	0
57	MG	YA	3010	1/1	0.97	0.13	3,3,3,3	0
57	MG	RA	3115	1/1	0.97	0.11	16,16,16,16	0
57	MG	QA	1605	1/1	0.97	0.22	17,17,17,17	0
57	MG	XA	1607	1/1	0.97	0.07	61,61,61,61	0
57	MG	RA	3011	1/1	0.97	0.18	11,11,11,11	0
57	MG	XA	1609	1/1	0.97	0.08	1,1,1,1	0
57	MG	RA	3063	1/1	0.97	0.33	1,1,1,1	0
57	MG	RA	3078	1/1	0.97	0.08	11,11,11,11	0
57	MG	YA	3199	1/1	0.97	0.11	21,21,21,21	0
57	MG	YA	3075	1/1	0.97	0.20	4,4,4,4	0
57	MG	YA	3140	1/1	0.97	0.08	47,47,47,47	0
57	MG	RA	3064	1/1	0.97	0.30	3,3,3,3	0
57	MG	YA	3019	1/1	0.97	0.18	3,3,3,3	0
57	MG	RA	3145	1/1	0.97	0.14	14,14,14,14	0
57	MG	XA	1648	1/1	0.97	0.12	21,21,21,21	0
57	MG	YA	3145	1/1	0.97	0.08	2,2,2,2	0
57	MG	XA	1692	1/1	0.97	0.12	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3146	1/1	0.97	0.14	38,38,38,38	0
57	MG	YA	3082	1/1	0.97	0.15	7,7,7,7	0
57	MG	YA	3282	1/1	0.97	0.30	48,48,48,48	0
57	MG	YA	3211	1/1	0.97	0.05	13,13,13,13	0
57	MG	XA	1694	1/1	0.97	0.05	13,13,13,13	0
57	MG	QA	1609	1/1	0.97	0.05	4,4,4,4	0
57	MG	RA	3199	1/1	0.97	0.14	44,44,44,44	0
57	MG	YA	3030	1/1	0.97	0.19	7,7,7,7	0
57	MG	RA	3067	1/1	0.97	0.06	4,4,4,4	0
57	MG	XA	1698	1/1	0.97	0.09	24,24,24,24	0
57	MG	QA	1649	1/1	0.97	0.23	32,32,32,32	0
57	MG	YA	3291	1/1	0.97	0.21	28,28,28,28	0
57	MG	RA	3249	1/1	0.98	0.08	44,44,44,44	0
57	MG	XA	1672	1/1	0.98	0.10	10,10,10,10	0
57	MG	XA	1673	1/1	0.98	0.11	36,36,36,36	0
57	MG	RA	3178	1/1	0.98	0.09	25,25,25,25	0
57	MG	YA	3241	1/1	0.98	0.20	3,3,3,3	0
57	MG	XA	1640	1/1	0.98	0.10	24,24,24,24	0
57	MG	YA	3133	1/1	0.98	0.05	0,0,0,0	0
57	MG	YA	3245	1/1	0.98	0.20	11,11,11,11	0
57	MG	RA	3142	1/1	0.98	0.15	11,11,11,11	0
57	MG	XA	1610	1/1	0.98	0.11	13,13,13,13	0
57	MG	YA	3136	1/1	0.98	0.09	2,2,2,2	0
57	MG	RA	3108	1/1	0.98	0.06	7,7,7,7	0
57	MG	QA	1616	1/1	0.98	0.05	7,7,7,7	0
57	MG	RA	3182	1/1	0.98	0.08	37,37,37,37	0
57	MG	YA	3253	1/1	0.98	0.17	7,7,7,7	0
57	MG	QA	1606	1/1	0.98	0.08	3,3,3,3	0
57	MG	XA	1616	1/1	0.98	0.13	7,7,7,7	0
57	MG	QM	201	1/1	0.98	0.10	45,45,45,45	0
57	MG	YA	3090	1/1	0.98	0.10	22,22,22,22	0
57	MG	YA	3091	1/1	0.98	0.14	3,3,3,3	0
57	MG	QA	1607	1/1	0.98	0.04	22,22,22,22	0
57	MG	RA	3113	1/1	0.98	0.07	9,9,9,9	0
57	MG	RA	3024	1/1	0.98	0.12	4,4,4,4	0
57	MG	YA	3047	1/1	0.98	0.20	7,7,7,7	0
57	MG	YA	3265	1/1	0.98	0.09	23,23,23,23	0
57	MG	RA	3188	1/1	0.98	0.09	25,25,25,25	0
57	MG	QA	1638	1/1	0.98	0.10	38,38,38,38	0
57	MG	YA	3008	1/1	0.98	0.12	3,3,3,3	0
57	MG	RA	3169	1/1	0.98	0.10	21,21,21,21	0
57	MG	YA	3101	1/1	0.98	0.11	25,25,25,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3288	1/1	0.98	0.20	34,34,34,34	0
57	MG	RA	3097	1/1	0.98	0.19	10,10,10,10	0
57	MG	RA	3192	1/1	0.98	0.05	21,21,21,21	0
57	MG	QA	1634	1/1	0.98	0.06	24,24,24,24	0
57	MG	XA	1659	1/1	0.98	0.21	35,35,35,35	0
57	MG	YA	3057	1/1	0.98	0.20	4,4,4,4	0
57	MG	YA	3108	1/1	0.98	0.05	15,15,15,15	0
57	MG	RA	3035	1/1	0.98	0.17	3,3,3,3	0
57	MG	RA	3217	1/1	0.98	0.20	6,6,6,6	0
57	MG	XA	1662	1/1	0.98	0.03	20,20,20,20	0
57	MG	RA	3218	1/1	0.98	0.28	22,22,22,22	0
57	MG	XA	1700	1/1	0.98	0.07	54,54,54,54	0
57	MG	YA	3063	1/1	0.98	0.10	4,4,4,4	0
57	MG	YA	3115	1/1	0.98	0.24	29,29,29,29	0
57	MG	RA	3077	1/1	0.98	0.14	3,3,3,3	0
57	MG	YA	3021	1/1	0.98	0.08	3,3,3,3	0
57	MG	YA	3118	1/1	0.98	0.19	19,19,19,19	0
57	MG	YA	3066	1/1	0.98	0.30	29,29,29,29	0
57	MG	XA	1601	1/1	0.98	0.10	7,7,7,7	0
57	MG	RA	3102	1/1	0.98	0.04	3,3,3,3	0
57	MG	YA	3230	1/1	0.98	0.14	8,8,8,8	0
57	MG	XA	1704	1/1	0.98	0.14	47,47,47,47	0
57	MG	XA	1603	1/1	0.98	0.12	9,9,9,9	0
57	MG	RA	3038	1/1	0.98	0.12	7,7,7,7	0
57	MG	YA	3027	1/1	0.98	0.16	2,2,2,2	0
57	MG	RA	3079	1/1	0.98	0.09	16,16,16,16	0
59	ZN	QN	101	1/1	0.98	0.04	89,89,89,89	0
57	MG	RA	3106	1/1	0.98	0.10	16,16,16,16	0
57	MG	QA	1640	1/1	0.99	0.09	8,8,8,8	0
57	MG	RA	3012	1/1	0.99	0.09	2,2,2,2	0
57	MG	YA	3028	1/1	0.99	0.11	3,3,3,3	0
57	MG	RA	3220	1/1	0.99	0.18	29,29,29,29	0
57	MG	YA	3212	1/1	0.99	0.07	22,22,22,22	0
57	MG	YA	3186	1/1	0.99	0.04	8,8,8,8	0
57	MG	RA	3026	1/1	0.99	0.04	8,8,8,8	0
57	MG	RA	3049	1/1	0.99	0.06	3,3,3,3	0
57	MG	RA	3100	1/1	0.99	0.03	2,2,2,2	0
57	MG	YA	3244	1/1	0.99	0.17	6,6,6,6	0
57	MG	YA	3073	1/1	0.99	0.06	17,17,17,17	0
57	MG	YA	3096	1/1	0.99	0.04	13,13,13,13	0
57	MG	RA	3143	1/1	0.99	0.06	22,22,22,22	0
57	MG	RA	3006	1/1	0.99	0.20	7,7,7,7	0

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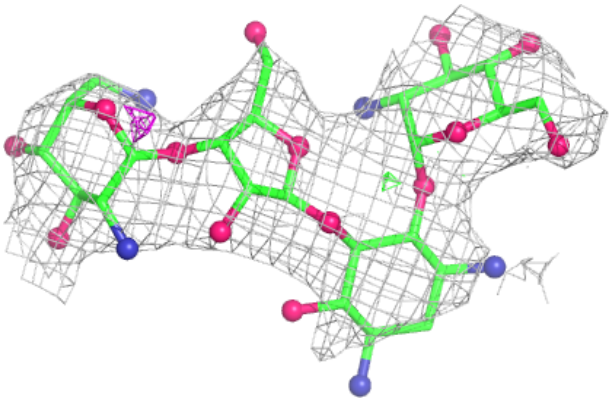
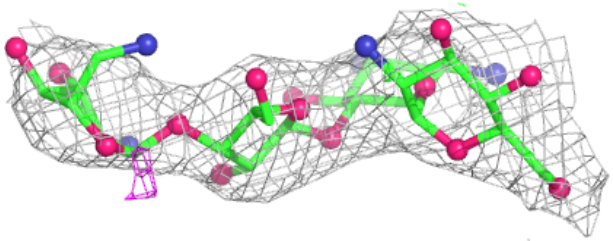
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
57	MG	RA	3017	1/1	0.99	0.09	2,2,2,2	0
57	MG	YA	3146	1/1	0.99	0.04	5,5,5,5	0
57	MG	YA	3251	1/1	0.99	0.30	36,36,36,36	0
57	MG	RA	3116	1/1	0.99	0.07	33,33,33,33	0
57	MG	RA	3103	1/1	0.99	0.04	9,9,9,9	0
57	MG	YA	3254	1/1	0.99	0.08	5,5,5,5	0
57	MG	RA	3018	1/1	0.99	0.05	9,9,9,9	0
57	MG	RA	3133	1/1	0.99	0.10	32,32,32,32	0
57	MG	XA	1612	1/1	0.99	0.03	2,2,2,2	0
57	MG	YA	3201	1/1	0.99	0.12	43,43,43,43	0
57	MG	RA	3036	1/1	0.99	0.12	2,2,2,2	0
57	MG	RA	3037	1/1	0.99	0.09	7,7,7,7	0
57	MG	RA	3107	1/1	0.99	0.05	18,18,18,18	0
57	MG	YA	3005	1/1	0.99	0.09	1,1,1,1	0
59	ZN	QD	301	1/1	0.99	0.15	37,37,37,37	0
57	MG	YA	3086	1/1	0.99	0.09	1,1,1,1	0
57	MG	RA	3065	1/1	0.99	0.12	3,3,3,3	0
59	ZN	XN	101	1/1	0.99	0.03	73,73,73,73	0
57	MG	RA	3051	1/1	1.00	0.08	3,3,3,3	0
57	MG	XA	1682	1/1	1.00	0.15	3,3,3,3	0
57	MG	YA	3255	1/1	1.00	0.07	17,17,17,17	0

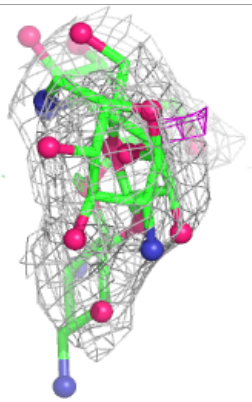
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PAR QA 1696:

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

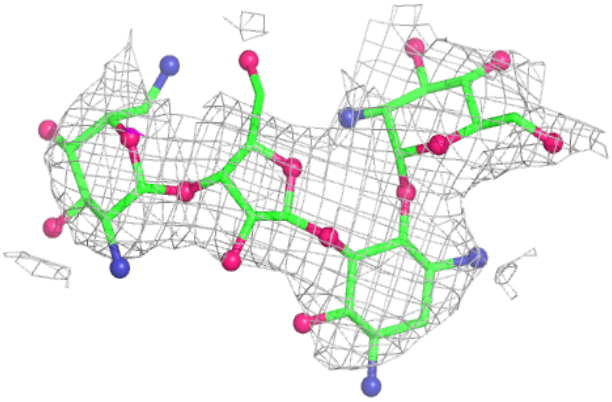


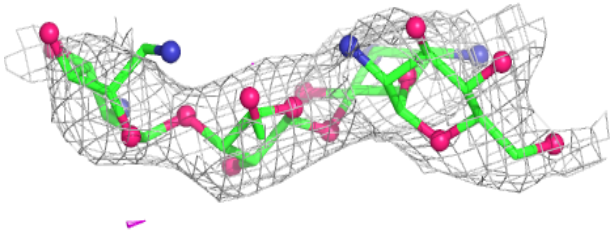


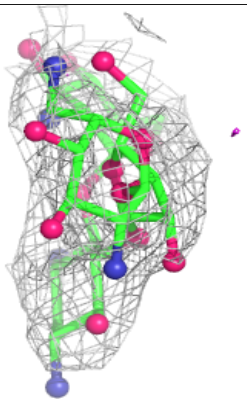


Electron density around PAR XA 1715:

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







6.5 Other polymers (i)

There are no such residues in this entry.