



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 14, 2016 – 07:23 PM EST

PDB ID : 5DGF  
Title : Complex of yeast 80S ribosome with hypusine-containing/non-modified eIF5A and/or a peptidyl-tRNA analog  
Authors : Melnikov, S.; Mailliot, J.; Shin, B.-S.; Rigger, L.; Yusupova, G.; Micura, R.; Dever, T.E.; Yusupov, M.  
Deposited on : 2015-08-27  
Resolution : 3.30 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20028442

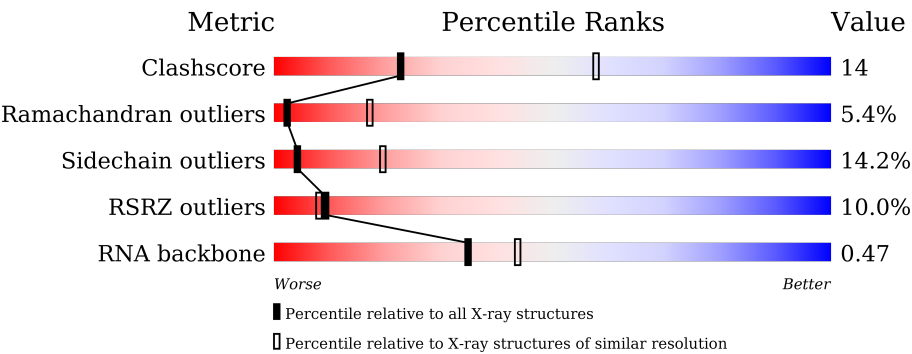
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)
RNA backbone	2183	1005 (3.82-2.78)

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

Mol	Chain	Length	Quality of chain
1	2	1800	<div><div>5%</div><div><div>38%</div><div>47%</div><div>14%</div><div>..</div></div></div>
1	6	1800	<div><div>3%</div><div><div>40%</div><div>45%</div><div>14%</div><div>.</div></div></div>
2	S0	251	<div><div>20%</div><div><div>27%</div><div>41%</div><div>14%</div><div>18%</div></div></div>
2	s0	251	<div><div>20%</div><div><div>67%</div><div>15%</div><div>18%</div></div></div>
3	S1	254	<div><div>30%</div><div><div>26%</div><div>47%</div><div>10%</div><div>16%</div></div></div>
3	s1	254	<div><div>17%</div><div><div>69%</div><div>15%</div><div>15%</div></div></div>

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Mol	Chain	Length	Quality of chain
4	S2	253	
4	s2	253	
5	S3	239	
5	s3	239	
6	S4	260	
6	s4	260	
7	S5	224	
7	s5	224	
8	S6	236	
8	s6	236	
9	S7	189	
9	s7	189	
10	S8	200	
10	s8	200	
11	S9	196	
11	s9	196	
12	C0	96	
12	c0	96	
13	C1	155	
13	c1	155	
14	C2	142	
14	c2	142	
15	C3	150	
15	c3	150	
16	C4	136	

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Mol	Chain	Length	Quality of chain
16	c4	136	
17	C5	141	
17	c5	141	
18	C6	142	
18	c6	142	
19	C7	136	
19	c7	136	
20	C8	145	
20	c8	145	
21	C9	143	
21	c9	143	
22	D0	120	
22	d0	120	
23	D1	87	
23	d1	87	
24	D2	129	
24	d2	129	
25	D3	144	
25	d3	144	
26	D4	134	
26	d4	134	
27	D5	107	
27	d5	107	
28	D6	97	
28	d6	97	

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Mol	Chain	Length	Quality of chain
29	D7	81	
29	d7	81	
30	D8	66	
30	d8	66	
31	D9	55	
31	d9	55	
32	E0	62	
32	e0	62	
33	E1	76	
33	e1	76	
34	SR	318	
34	sR	318	
35	SM	273	
35	sM	273	
36	1	3396	
36	5	3396	
37	3	121	
37	7	121	
38	4	158	
38	8	158	
39	L2	253	
39	l2	253	
40	L3	386	
40	l3	386	
41	L4	361	

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Mol	Chain	Length	Quality of chain
41	l4	361	
42	L5	296	
42	l5	296	
43	L6	175	
43	l6	175	
44	L7	243	
44	l7	243	
45	L8	255	
45	l8	255	
46	L9	191	
46	l9	191	
47	M0	220	
47	m0	220	
48	M1	173	
48	m1	173	
49	M3	198	
49	m3	198	
50	M4	137	
50	m4	137	
51	M5	203	
51	m5	203	
52	M6	198	
52	m6	198	
53	M7	183	
53	m7	183	

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Mol	Chain	Length	Quality of chain
54	M8	185	
54	m8	185	
55	M9	188	
55	m9	188	
56	N0	172	
56	n0	172	
57	N1	159	
57	n1	159	
58	N2	120	
58	n2	120	
59	N3	136	
59	n3	136	
60	N4	155	
60	n4	155	
61	N5	141	
61	n5	141	
62	N6	126	
62	n6	126	
63	N7	135	
63	n7	135	
64	N8	148	
64	n8	148	
65	N9	58	
65	n9	58	
66	O0	104	

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Mol	Chain	Length	Quality of chain
66	o0	104	
67	O1	112	
67	o1	112	
68	O2	129	
68	o2	129	
69	O3	106	
69	o3	106	
70	O4	120	
70	o4	120	
71	O5	119	
71	o5	119	
72	O6	99	
72	o6	99	
73	O7	87	
73	o7	87	
74	O8	77	
74	o8	77	
75	O9	50	
75	o9	50	
76	Q0	52	
76	q0	52	
77	Q1	25	
77	q1	25	
78	Q2	105	
78	q2	105	

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Mol	Chain	Length	Quality of chain
79	Q3	91	
79	q3	91	
80	m2	165	
81	p0	311	
82	p1	47	
82	p2	47	
83	f	157	
84	B	5	
84	C	5	

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
86	SPS	B	3401	-	-	-	X
87	MG	5	3401	-	-	-	X
87	MG	5	3404	-	-	-	X
87	MG	5	3405	-	-	-	X

## 2 Entry composition

There are 89 unique types of molecules in this entry. The entry contains 404292 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 18S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	2	1781	Total	C	N	O	P	0	1	0
			37970	16975	6720	12493	1782			
1	6	1795	Total	C	N	O	P	0	1	0
			38260	17105	6763	12596	1796			

- Molecule 2 is a protein called 40S ribosomal protein S0-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S0	206	Total	C	N	O	S	0	0	0
			1577	1014	278	283	2			
2	s0	206	Total	C	N	O	S	0	0	0
			1583	1017	281	283	2			

- Molecule 3 is a protein called 40S ribosomal protein S1-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	S1	214	Total	C	N	O	S	0	0	0
			1709	1084	310	311	4			
3	s1	216	Total	C	N	O	S	0	0	0
			1722	1091	312	315	4			

- Molecule 4 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	S2	217	Total	C	N	O	S	0	0	0
			1635	1047	289	297	2			
4	s2	217	Total	C	N	O	S	0	0	0
			1635	1047	289	297	2			

- Molecule 5 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	S3	223	Total	C	N	O	S	0	0	0
			1734	1101	313	314	6			
5	s3	223	Total	C	N	O	S	0	0	0
			1734	1101	313	314	6			

- Molecule 6 is a protein called 40S ribosomal protein S4-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	S4	260	Total	C	N	O	S	0	0	0
			2068	1316	389	360	3			
6	s4	260	Total	C	N	O	S	0	0	0
			2068	1316	389	360	3			

- Molecule 7 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	S5	206	Total	C	N	O	S	0	0	0
			1609	1007	300	299	3			
7	s5	206	Total	C	N	O	S	0	0	0
			1609	1007	300	299	3			

- Molecule 8 is a protein called 40S ribosomal protein S6-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	S6	226	Total	C	N	O	S	0	0	0
			1799	1129	346	321	3			
8	s6	218	Total	C	N	O	S	0	0	0
			1755	1102	337	313	3			

- Molecule 9 is a protein called 40S ribosomal protein S7-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	S7	184	Total	C	N	O	S	0	0	0
			1481	951	265	265				
9	s7	186	Total	C	N	O	S	0	0	0
			1491	957	267	267				

- Molecule 10 is a protein called 40S ribosomal protein S8-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	S8	188	Total	C	N	O	S	0	0	0
			1489	925	298	264	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	s8	188	Total	C	N	O	S	0	0	0
			1489	925	298	264	2			

- Molecule 11 is a protein called 40S ribosomal protein S9-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	S9	185	Total	C	N	O	S	0	0	0
			1494	943	289	261	1			
11	s9	185	Total	C	N	O	S	0	0	0
			1494	943	289	261	1			

- Molecule 12 is a protein called 40S ribosomal protein S10-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	C0	96	Total	C	N	O	S	0	0	0
			773	500	126	145	2			
12	c0	96	Total	C	N	O	S	0	0	0
			762	491	125	144	2			

- Molecule 13 is a protein called 40S ribosomal protein S11-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	C1	155	Total	C	N	O	S	0	0	0
			1214	775	230	206	3			
13	c1	146	Total	C	N	O	S	0	0	0
			1168	747	221	197	3			

- Molecule 14 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	C2	124	Total	C	N	O	S	0	0	0
			892	562	156	172	2			
14	c2	124	Total	C	N	O	S	0	0	0
			892	562	156	172	2			

- Molecule 15 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	C3	150	Total	C	N	O	S	0	0	0
			1192	759	224	207	2			
15	c3	150	Total	C	N	O	S	0	0	0
			1192	759	224	207	2			



- Molecule 16 is a protein called 40S ribosomal protein S14-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	C4	127	Total	C	N	O	S	0	0	0
			891	545	182	163	1			
16	c4	128	Total	C	N	O	S	0	0	0
			949	582	188	176	3			

- Molecule 17 is a protein called 40S ribosomal protein S15 (uS19).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	C5	124	Total	C	N	O	S	0	0	0
			977	622	182	166	7			
17	c5	135	Total	C	N	O	S	0	0	0
			1039	658	196	178	7			

- Molecule 18 is a protein called 40S ribosomal protein S16-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	C6	141	Total	C	N	O		0	0	0
			1105	708	203	194				
18	c6	142	Total	C	N	O		0	0	0
			1111	711	204	196				

- Molecule 19 is a protein called 40S ribosomal protein S17-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	C7	120	Total	C	N	O	S	0	0	0
			926	577	177	170	2			
19	c7	117	Total	C	N	O	S	0	0	0
			906	563	174	167	2			

- Molecule 20 is a protein called 40S ribosomal protein S18-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	C8	145	Total	C	N	O	S	0	0	0
			1192	743	237	210	2			
20	c8	145	Total	C	N	O	S	0	0	0
			1192	743	237	210	2			

- Molecule 21 is a protein called 40S ribosomal protein S19-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	C9	143	Total	C	N	O	S	0	0	0
			1112	694	208	208	2			
21	c9	143	Total	C	N	O	S	0	0	0
			1112	694	208	208	2			

- Molecule 22 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	D0	107	Total	C	N	O	S	0	0	0
			855	539	156	159	1			
22	d0	110	Total	C	N	O	S	0	0	0
			882	554	161	166	1			

- Molecule 23 is a protein called 40S ribosomal protein S21-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	D1	87	Total	C	N	O	S	0	0	0
			684	420	125	137	2			
23	d1	87	Total	C	N	O	S	0	0	0
			684	420	125	137	2			

- Molecule 24 is a protein called 40S ribosomal protein S22-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	D2	129	Total	C	N	O	S	0	0	0
			1021	650	188	180	3			
24	d2	129	Total	C	N	O	S	0	0	0
			1021	650	188	180	3			

- Molecule 25 is a protein called 40S ribosomal protein S23-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	D3	144	Total	C	N	O	S	0	0	0
			1121	708	220	191	2			
25	d3	144	Total	C	N	O	S	0	0	0
			1121	708	220	191	2			

- Molecule 26 is a protein called 40S ribosomal protein S24-A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
26	D4	134	Total	C	N	O	0	0	0
			1073	676	208	189			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
26	d4	134	Total	C	N	O	0	0	0
			1073	676	208	189			

- Molecule 27 is a protein called 40S ribosomal protein S25-A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
27	D5	70	Total	C	N	O	0	0	0
			563	360	104	99			
27	d5	69	Total	C	N	O	0	0	0
			558	357	103	98			

- Molecule 28 is a protein called 40S ribosomal protein S26-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	D6	97	Total	C	N	O	S	0	0	0
			769	475	160	129	5			
28	d6	97	Total	C	N	O	S	0	0	0
			769	475	160	129	5			

- Molecule 29 is a protein called 40S ribosomal protein S27-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	D7	81	Total	C	N	O	S	0	0	0
			610	382	110	113	5			
29	d7	81	Total	C	N	O	S	0	0	0
			610	382	110	113	5			

- Molecule 30 is a protein called 40S ribosomal protein S28-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	D8	63	Total	C	N	O	S	0	0	0
			497	306	99	91	1			
30	d8	63	Total	C	N	O	S	0	0	0
			497	306	99	91	1			

- Molecule 31 is a protein called 40S ribosomal protein S29-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	D9	53	Total	C	N	O	S	0	0	0
			442	274	92	72	4			
31	d9	53	Total	C	N	O	S	0	0	0
			442	274	92	72	4			

- Molecule 32 is a protein called 40S ribosomal protein S30-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	E0	60	Total	C	N	O	S	0	0	0
			475	299	98	77	1			
32	e0	62	Total	C	N	O	S	0	0	0
			491	309	101	80	1			

- Molecule 33 is a protein called Ubiquitin-40S ribosomal protein S31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	E1	71	Total	C	N	O	S	0	0	0
			566	362	106	94	4			
33	e1	76	Total	C	N	O	S	0	0	0
			608	388	117	99	4			

- Molecule 34 is a protein called Guanine nucleotide-binding protein subunit beta-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	SR	318	Total	C	N	O	S	0	0	0
			2437	1541	418	470	8			
34	sR	318	Total	C	N	O	S	0	0	0
			2442	1544	418	472	8			

- Molecule 35 is a protein called Suppressor protein STM1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	SM	159	Total	C	N	O		0	0	0
			1105	653	221	231				
35	sM	104	Total	C	N	O		0	0	0
			680	403	140	137				

- Molecule 36 is a RNA chain called 25S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	1	3149	Total	C	N	O	P	0	0	0
			67355	30086	12142	21978	3149			
36	5	3169	Total	C	N	O	P	0	0	0
			67780	30276	12216	22120	3168			

- Molecule 37 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	3	121	Total	C	N	O	P	0	0	0
			2579	1152	461	845	121			
37	7	121	Total	C	N	O	P	0	0	0
			2579	1152	461	845	121			

- Molecule 38 is a RNA chain called 5.8S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	4	158	Total	C	N	O	P	0	0	0
			3353	1500	586	1109	158			
38	8	158	Total	C	N	O	P	0	0	0
			3353	1500	586	1109	158			

- Molecule 39 is a protein called 60S ribosomal protein L2-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	L2	252	Total	C	N	O	S	0	0	0
			1914	1191	388	334	1			
39	12	252	Total	C	N	O	S	0	0	0
			1912	1190	388	333	1			

- Molecule 40 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	L3	386	Total	C	N	O	S	0	0	0
			3075	1950	584	533	8			
40	13	386	Total	C	N	O	S	0	0	0
			3075	1950	584	533	8			

- Molecule 41 is a protein called 60S ribosomal protein L4-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	L4	361	Total	C	N	O	S	0	0	0
			2748	1729	522	494	3			
41	14	361	Total	C	N	O	S	0	0	0
			2748	1729	522	494	3			

- Molecule 42 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	L5	296	Total	C	N	O	S	0	0	0
			2375	1501	414	458	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	15	294	Total	C	N	O	S	0	0	0
			2359	1489	412	456	2			

- Molecule 43 is a protein called 60S ribosomal protein L6-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	L6	156	Total	C	N	O	S	0	0	0
			1239	800	222	216	1			
43	16	157	Total	C	N	O	S	0	0	0
			1248	806	224	217	1			

- Molecule 44 is a protein called 60S ribosomal protein L7-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	L7	222	Total	C	N	O	S	0	0	0
			1784	1151	324	308	1			
44	17	223	Total	C	N	O	S	0	0	0
			1791	1155	325	310	1			

- Molecule 45 is a protein called 60S ribosomal protein L8-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	L8	233	Total	C	N	O	S	0	0	0
			1804	1151	323	327	3			
45	18	231	Total	C	N	O	S	0	0	0
			1764	1131	316	314	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L8	119	ALA	GLY	conflict	UNP P17076
18	119	ALA	GLY	conflict	UNP P17076

- Molecule 46 is a protein called 60S ribosomal protein L9-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	L9	191	Total	C	N	O	S	0	0	0
			1518	963	274	277	4			
46	19	191	Total	C	N	O	S	0	0	0
			1518	963	274	277	4			

- Molecule 47 is a protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	M0	211	Total	C	N	O	S	0	0	0
			1705	1083	322	294	6			
47	m0	213	Total	C	N	O	S	0	0	0
			1722	1094	325	297	6			

- Molecule 48 is a protein called 60S ribosomal protein L11-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	M1	169	Total	C	N	O	S	0	0	0
			1353	847	253	249	4			
48	m1	169	Total	C	N	O	S	0	0	0
			1353	847	253	249	4			

- Molecule 49 is a protein called 60S ribosomal protein L13-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	M3	193	Total	C	N	O		0	0	0
			1543	962	315	266				
49	m3	194	Total	C	N	O		0	0	0
			1548	965	316	267				

- Molecule 50 is a protein called 60S ribosomal protein L14-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	M4	136	Total	C	N	O	S	0	0	0
			1053	675	199	177	2			
50	m4	137	Total	C	N	O	S	0	0	0
			1059	678	200	179	2			

- Molecule 51 is a protein called 60S ribosomal protein L15-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	M5	203	Total	C	N	O	S	0	0	0
			1720	1077	361	281	1			
51	m5	203	Total	C	N	O	S	0	0	0
			1720	1077	361	281	1			

- Molecule 52 is a protein called 60S ribosomal protein L16-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	M6	197	Total	C	N	O	S	0	0	0
			1555	1003	289	262	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	m6	197	Total	C	N	O	S	0	0	0
			1555	1003	289	262	1			

- Molecule 53 is a protein called 60S ribosomal protein L17-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	M7	183	Total	C	N	O	S	0	0	0
			1420	882	281	257				
53	m7	155	Total	C	N	O	S	0	0	0
			1227	764	238	225				

- Molecule 54 is a protein called 60S ribosomal protein L18-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	M8	185	Total	C	N	O	S	0	0	0
			1441	908	290	241	2			
54	m8	185	Total	C	N	O	S	0	0	0
			1441	908	290	241	2			

- Molecule 55 is a protein called 60S ribosomal protein L19-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	M9	188	Total	C	N	O	S	0	0	0
			1521	935	326	260				
55	m9	188	Total	C	N	O	S	0	0	0
			1521	935	326	260				

- Molecule 56 is a protein called 60S ribosomal protein L20-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
56	N0	172	Total	C	N	O	S	0	0	0
			1445	930	267	244	4			
56	n0	172	Total	C	N	O	S	0	0	0
			1445	930	267	244	4			

- Molecule 57 is a protein called 60S ribosomal protein L21-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
57	N1	159	Total	C	N	O	S	0	0	0
			1276	805	246	221	4			
57	n1	159	Total	C	N	O	S	0	0	0
			1276	805	246	221	4			



- Molecule 58 is a protein called 60S ribosomal protein L22-A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
58	N2	100	Total	C	N	O	0	0	0
			796	516	131	149			
58	n2	98	Total	C	N	O	0	0	0
			778	505	127	146			

- Molecule 59 is a protein called 60S ribosomal protein L23-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
59	N3	136	Total	C	N	O	S	0	0	0
			1003	628	189	179	7			
59	n3	136	Total	C	N	O	S	0	0	0
			1003	628	189	179	7			

- Molecule 60 is a protein called 60S ribosomal protein L24-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
60	N4	98	Total	C	N	O	S	0	0	0
			699	443	137	118	1			
60	n4	135	Total	C	N	O	S	0	0	0
			1038	651	206	180	1			

- Molecule 61 is a protein called 60S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
61	N5	121	Total	C	N	O	S	0	0	0
			964	620	169	173	2			
61	n5	120	Total	C	N	O	S	0	0	0
			959	617	168	172	2			

- Molecule 62 is a protein called 60S ribosomal protein L26-A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
62	N6	126	Total	C	N	O	0	0	0
			993	625	192	176			
62	n6	126	Total	C	N	O	0	0	0
			993	625	192	176			

- Molecule 63 is a protein called 60S ribosomal protein L27-A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
63	N7	135	Total	C	N	O	0	0	0
			1092	710	202	180			
63	n7	135	Total	C	N	O	0	0	0
			1092	710	202	180			

- Molecule 64 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
64	N8	148	Total	C	N	O	S	0	0	0
			1173	749	231	190	3			
64	n8	148	Total	C	N	O	S	0	0	0
			1173	749	231	190	3			

- Molecule 65 is a protein called 60S ribosomal protein L29.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
65	N9	58	Total	C	N	O	0	0	0
			462	289	100	73			
65	n9	58	Total	C	N	O	0	0	0
			462	289	100	73			

- Molecule 66 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
66	O0	97	Total	C	N	O	S	0	0	0
			743	479	124	139	1			
66	o0	100	Total	C	N	O	S	0	0	0
			767	492	128	146	1			

- Molecule 67 is a protein called 60S ribosomal protein L31-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
67	O1	109	Total	C	N	O	S	0	0	0
			876	556	167	152	1			
67	o1	109	Total	C	N	O	S	0	0	0
			883	559	167	156	1			

- Molecule 68 is a protein called 60S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
68	O2	127	Total	C	N	O	S	0	0	0
			1020	647	205	167	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
68	o2	127	Total	C	N	O	S	0	0	0
			1020	647	205	167	1			

- Molecule 69 is a protein called 60S ribosomal protein L33-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
69	O3	106	Total	C	N	O	S	0	0	0
			850	540	165	144	1			
69	o3	106	Total	C	N	O	S	0	0	0
			850	540	165	144	1			

- Molecule 70 is a protein called 60S ribosomal protein L34-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
70	O4	112	Total	C	N	O	S	0	0	0
			880	545	179	152	4			
70	o4	112	Total	C	N	O	S	0	0	0
			880	545	179	152	4			

- Molecule 71 is a protein called 60S ribosomal protein L35-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
71	O5	119	Total	C	N	O	S	0	0	0
			969	615	186	167	1			
71	o5	119	Total	C	N	O	S	0	0	0
			965	612	185	167	1			

- Molecule 72 is a protein called 60S ribosomal protein L36-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
72	O6	99	Total	C	N	O	S	0	0	0
			771	481	156	132	2			
72	o6	99	Total	C	N	O	S	0	0	0
			770	481	156	131	2			

- Molecule 73 is a protein called 60S ribosomal protein L37-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
73	O7	87	Total	C	N	O	S	0	0	0
			681	414	148	114	5			
73	o7	87	Total	C	N	O	S	0	0	0
			681	414	148	114	5			

- Molecule 74 is a protein called 60S ribosomal protein L38.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
74	O8	77	Total	C	N	O	0	0	0
			612	391	115	106			
74	o8	77	Total	C	N	O	0	0	0
			608	388	114	106			

- Molecule 75 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
75	O9	50	Total	C	N	O	S	0	0	0
			436	272	97	65	2			
75	o9	50	Total	C	N	O	S	0	0	0
			436	272	97	65	2			

- Molecule 76 is a protein called Ubiquitin-60S ribosomal protein L40.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
76	Q0	52	Total	C	N	O	S	0	0	0
			417	259	86	67	5			
76	q0	52	Total	C	N	O	S	0	0	0
			417	259	86	67	5			

- Molecule 77 is a protein called 60S ribosomal protein L41-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
77	Q1	25	Total	C	N	O	S	0	0	0
			233	142	63	27	1			
77	q1	25	Total	C	N	O	S	0	0	0
			233	142	63	27	1			

- Molecule 78 is a protein called 60S ribosomal protein L42-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
78	Q2	105	Total	C	N	O	S	0	0	0
			847	534	170	138	5			
78	q2	105	Total	C	N	O	S	0	0	0
			847	534	170	138	5			

- Molecule 79 is a protein called 60S ribosomal protein L43-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
79	Q3	91	Total	C	N	O	S	0	0	0
			694	429	138	121	6			
79	q3	91	Total	C	N	O	S	0	0	0
			694	429	138	121	6			

- Molecule 80 is a protein called 60S ribosomal protein L12-A (uL11).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
80	m2	150	Total	C	N	O		0	0	0
			750	450	150	150				

- Molecule 81 is a protein called 60S acidic ribosomal protein P0.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
81	p0	143	Total	C	N	O	S	0	0	0
			1077	687	192	195	3			

- Molecule 82 is a protein called 60S ribosomal protein P1 alpha/P2 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
82	p1	47	Total	C	N	O		0	0	0
			235	141	47	47				
82	p2	46	Total	C	N	O		0	0	0
			230	138	46	46				

- Molecule 83 is a protein called Eukaryotic translation initiation factor 5A-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
83	f	148	Total	C	N	O	S	0	0	0
			1116	692	188	227	9			

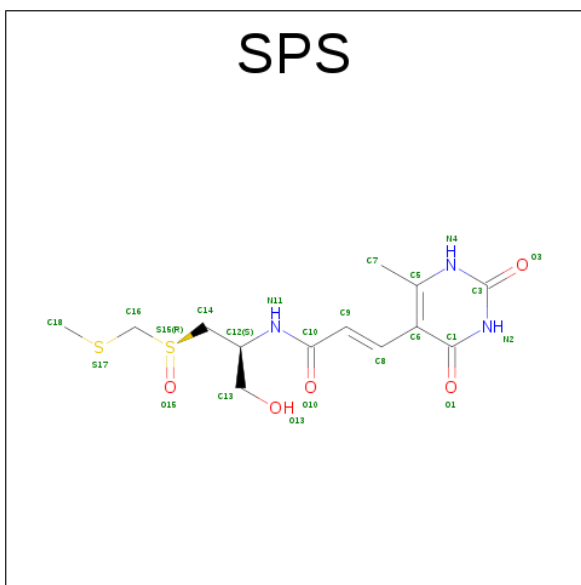
- Molecule 84 is a protein called CH-CH-8AN-PRO-PRO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
84	B	5	Total	C	N	O	P	0	0	0
			76	38	14	21	3			
84	C	5	Total	C	N	O	P	0	0	0
			73	38	14	19	2			

- Molecule 85 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
85	q0	1	Total Zn 1 1	0	0
85	D6	1	Total Zn 1 1	0	0
85	Q2	1	Total Zn 1 1	0	0
85	e1	1	Total Zn 1 1	0	0
85	Q3	1	Total Zn 1 1	0	0
85	D9	1	Total Zn 1 1	0	0
85	E1	1	Total Zn 1 1	0	0
85	Q0	1	Total Zn 1 1	0	0
85	d7	1	Total Zn 1 1	0	0
85	q3	1	Total Zn 1 1	0	0
85	d9	1	Total Zn 1 1	0	0
85	D7	1	Total Zn 1 1	0	0
85	d6	1	Total Zn 1 1	0	0
85	o7	1	Total Zn 1 1	0	0
85	O7	1	Total Zn 1 1	0	0
85	q2	1	Total Zn 1 1	0	0

- Molecule 86 is SPARSOMYCIN (three-letter code: SPS) (formula:  $C_{13}H_{19}N_3O_5S_2$ ).

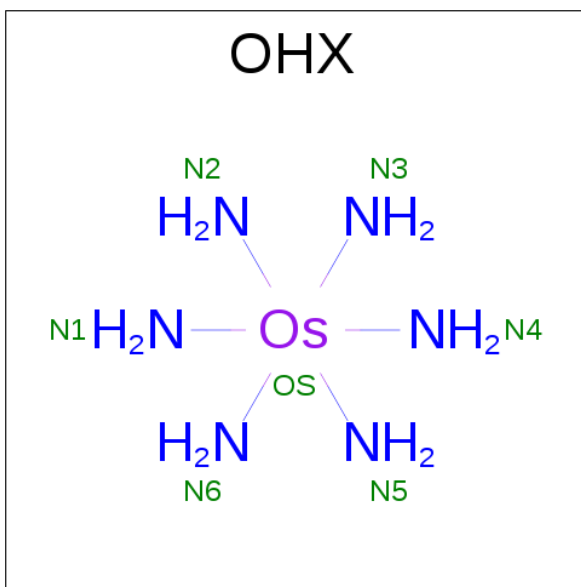


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
86	1	1	Total	C	N	O	S	0	0
			23	13	3	5	2		
86	B	1	Total	C	N	O	S	0	0
			23	13	3	5	2		

- Molecule 87 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
87	B	1	Total	Mg	0	0
			1	1		
87	1	3	Total	Mg	0	0
			3	3		
87	C	1	Total	Mg	0	0
			1	1		
87	f	2	Total	Mg	0	0
			2	2		
87	5	5	Total	Mg	0	0
			5	5		

- Molecule 88 is osmium (III) hexammine (three-letter code: OHX) (formula: H<sub>12</sub>N<sub>6</sub>Os).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
88	1	1	Total	N	Os	0	0
			7	6	1		
88	1	1	Total	N	Os	0	0
			7	6	1		
88	5	1	Total	N	Os	0	0
			7	6	1		

- Molecule 89 is water.

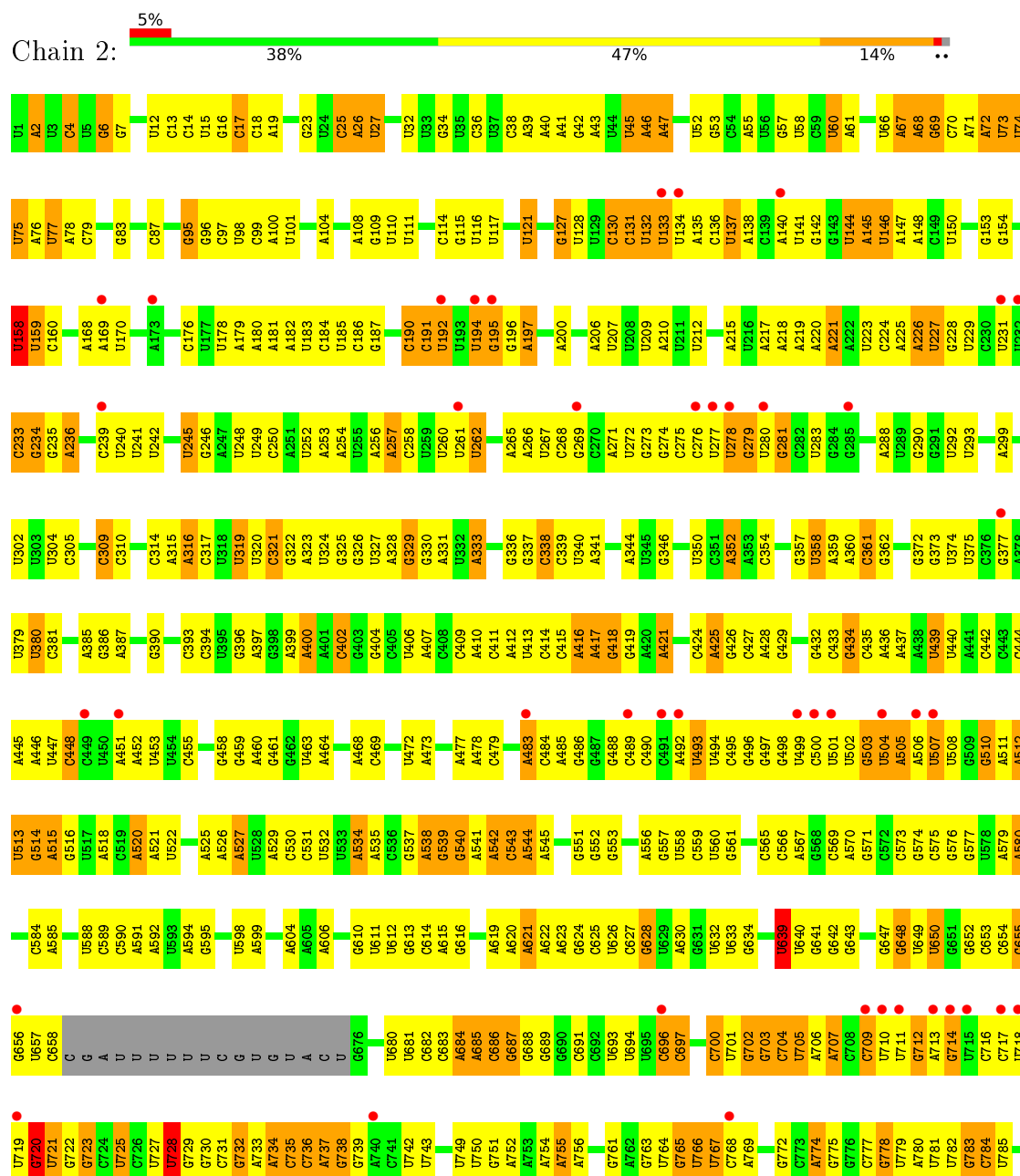
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
89	5	2	Total	O	0	0
			2	2		
89	f	10	Total	O	0	0
			10	10		

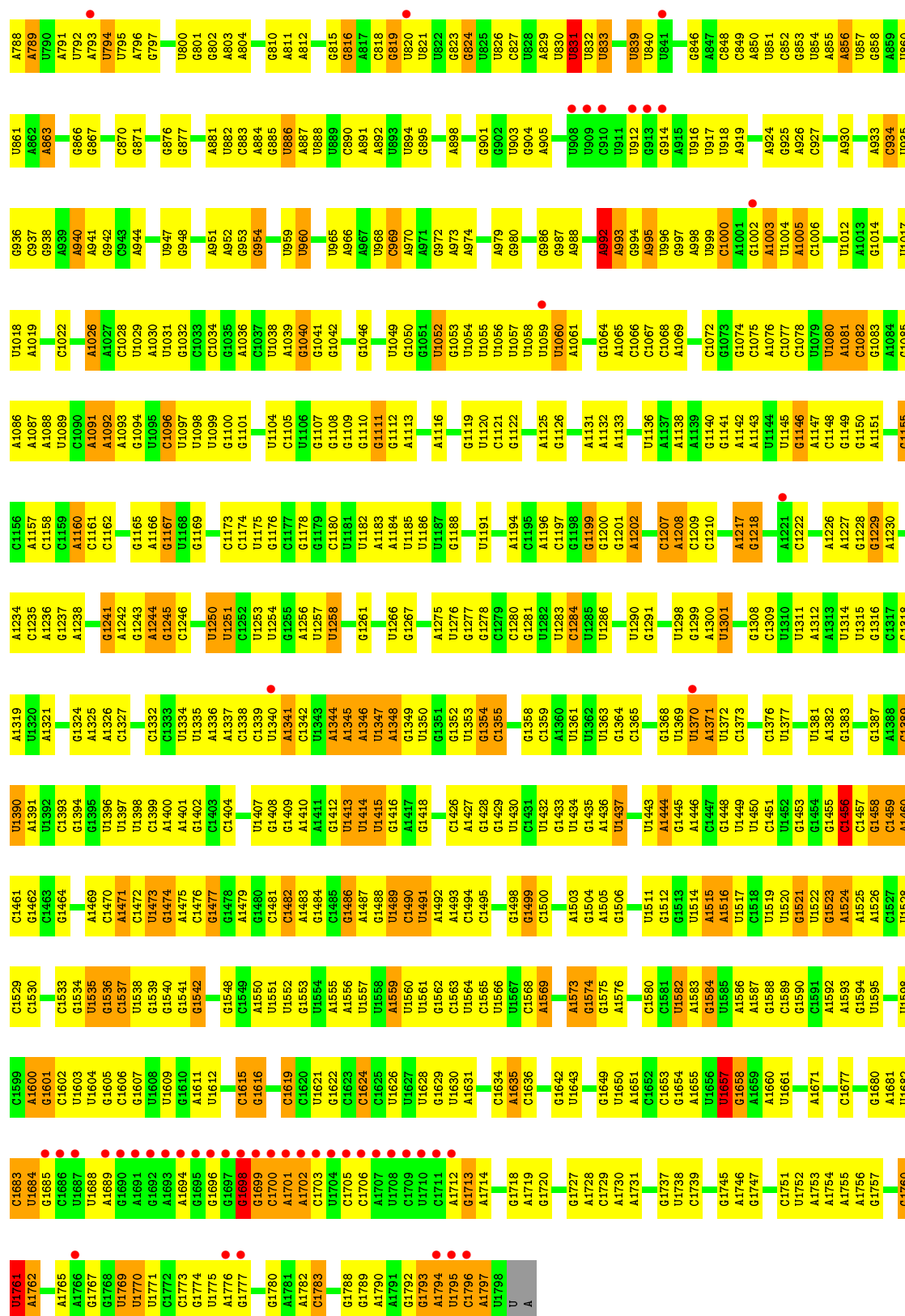


### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: 18S ribosomal RNA



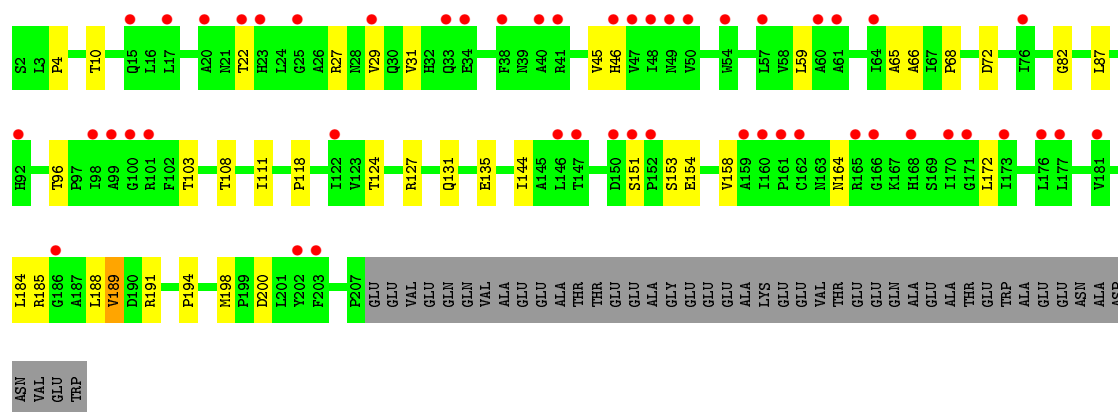


● Molecule 1: 18S ribosomal RNA

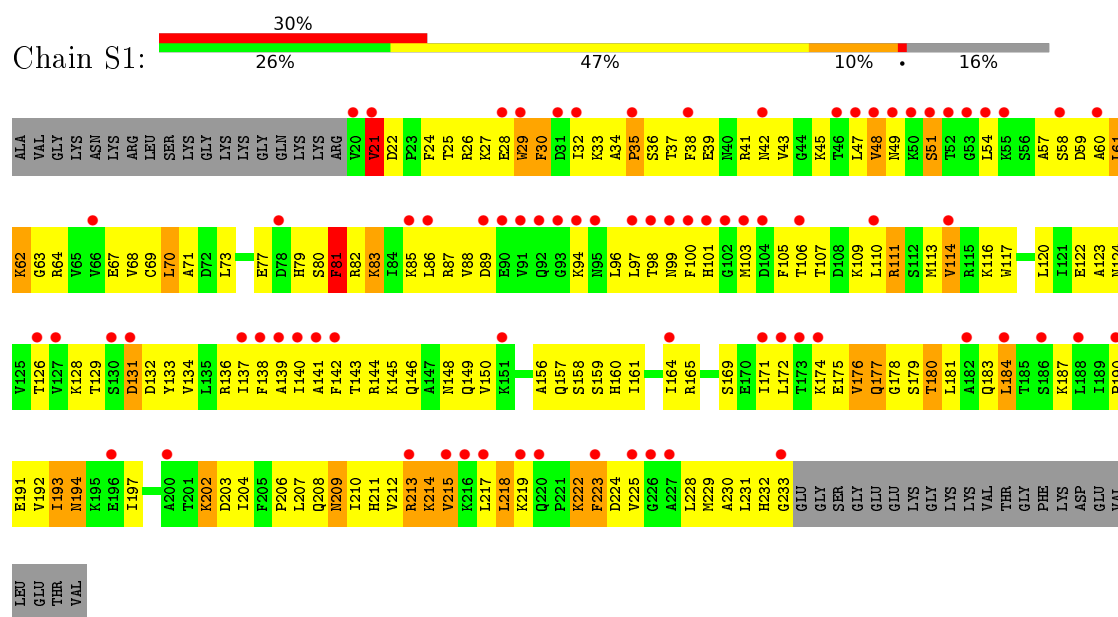


A1133	A1137	A1138	A1139	G1140	A1142	A1143	G1150	A1151	G1152	G1153	G1154	G1155	C1158	C1159	C1160	C1161	G1162	A1163	G1164	G1165	G1166	G1167	U1175	G1176	G1177	A1183	A1184	U1185	U1186	A1189	C1190	U1191	A1194	C1195	C1196	C1197	G1198	G1199	G1200	G1201	G1202	G1203	G1204	G1205	G1206	G1207	G1208	G1209	G1210	G1211	G1212	G1213	G1214	G1215	G1216																																																																																																																																																																																																																																																																																																																																																																																																											
U1063	A1064	A1065	C1066	C1067	C1068	A1069	C1070	U1071	U1072	G1073	G1074	C1075	A1076	C1077	C1078	C1079	C1080	A1081	C1082	G1083	A1084	A1085	A1086	A1087	A1088	A1089	A1090	A1091	A1092	U1095	C1096	U1097	U1098	U1099	G1100	G1101	G1102	U1103	G1107	G1108	G1109	G1110	G1111	G1112	A1113	G1114	G1115	A1116	U1117	G1118	G1119	G1120	G1121	G1122	G1123	G1124	G1125	G1126	G1127	G1130																																																																																																																																																																																																																																																																																																																																																																																																						
C897	A898	U899	A900	A901	G902	A903	A904	A905	U906	G907	G908	G909	G910	G911	G912	G913	G914	U917	U918	U919	U920	U921	U922	U923	U924	U925	U926	U927	U928	U929	U930	U931	U932	U933	U934	U935	U936	U937	U938	U939	U940	U941	U942	U943	U944	U945	U946	U947	U948	U949	U950	U951	U952	U953	U954	U955	U956	U957	U958	U959	U960	U961	U962	U963	U964	U965	U966	U967	U968	U969	U970	U971	U972	U973	U974	U975	U976	U977	U978	U979	U980	U981	U982	U983	U984	U985	U986	U987	U988	U989	U990	U991	U992	U993	U994	U995	U996	U997	U998	U999	U1000	U1001	U1002	U1003	U1004	U1005	U1006	U1007	U1008	U1009	U1010	U1011	U1012	U1013	U1014	U1015	U1016	U1017	U1018	U1019	U1020	U1021	U1022	U1023	U1024	U1025	U1026	U1027	U1028	U1029	U1030	U1031	U1032	U1033	U1034	U1035	U1036	U1037	U1038	U1039	U1040	U1041	U1042	U1043	U1044	U1045	U1046	U1047	U1048	U1049	U1050	U1051	U1052	U1053	U1054	U1055	U1056	U1057	U1058	U1059	U1060																																																																																																																																																																																																																																																																																																	
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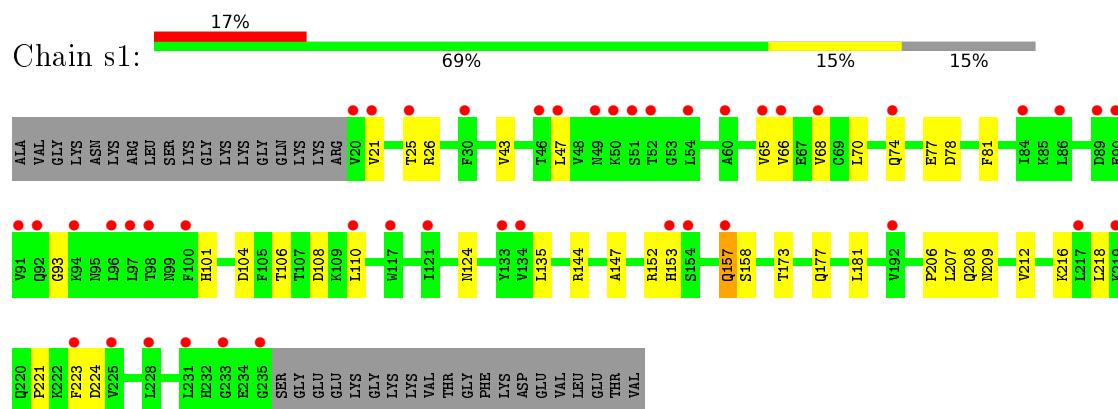




• Molecule 3: 40S ribosomal protein S1-A

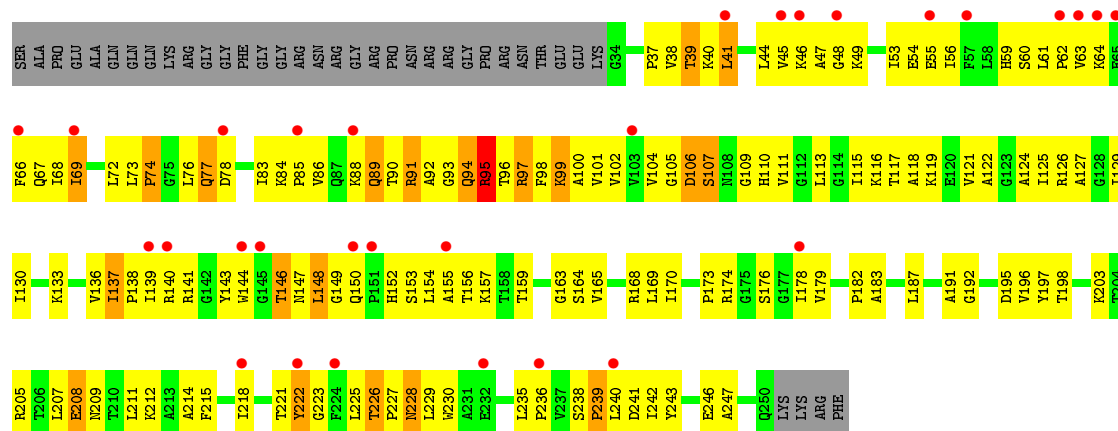


• Molecule 3: 40S ribosomal protein S1-A

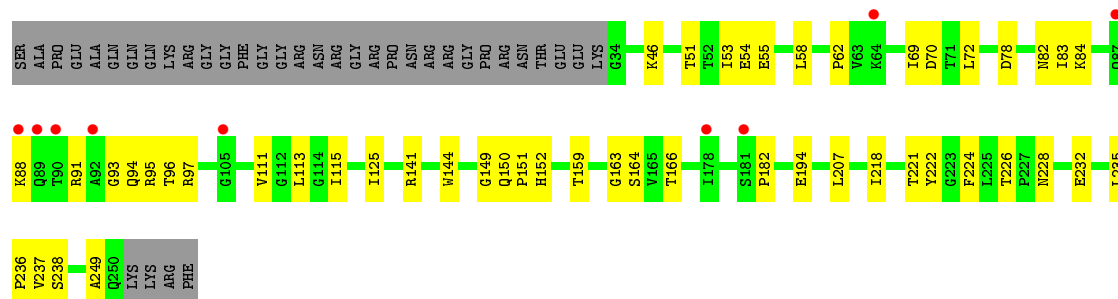


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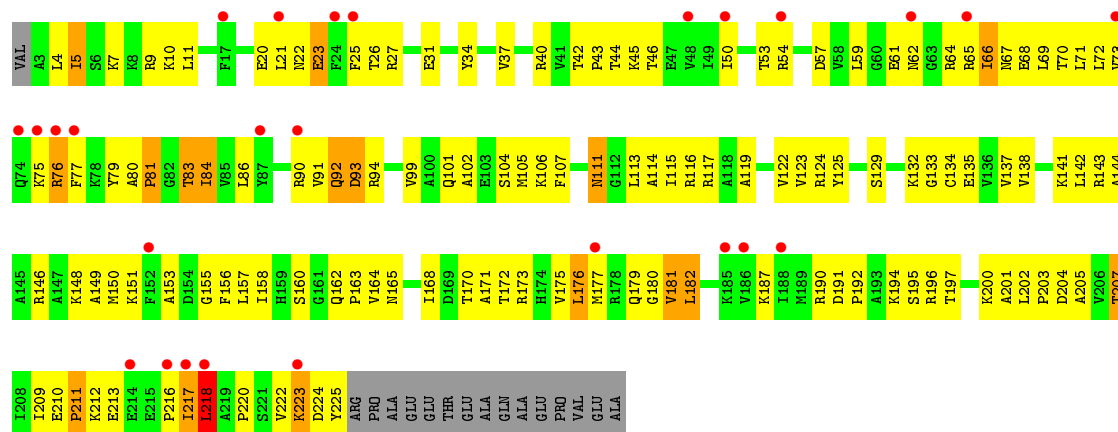




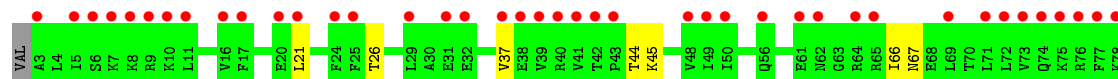
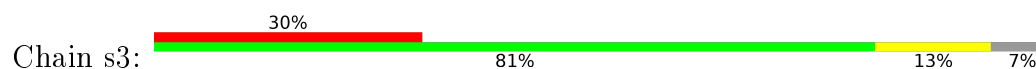
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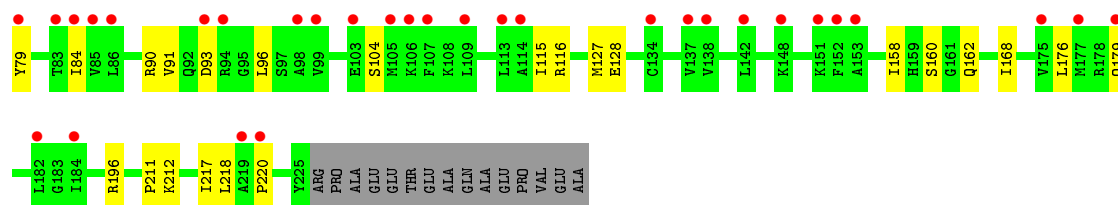


• Molecule 5: 40S ribosomal protein S3

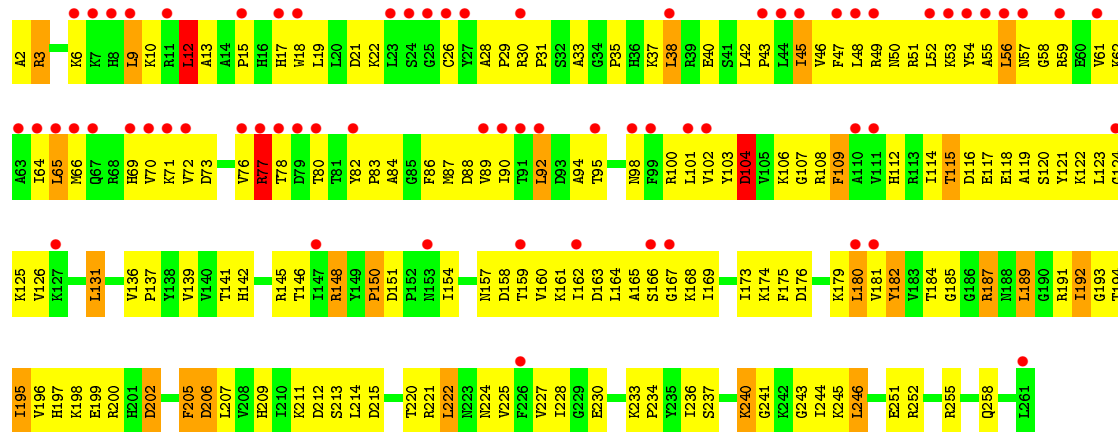


• Molecule 5: 40S ribosomal protein S3

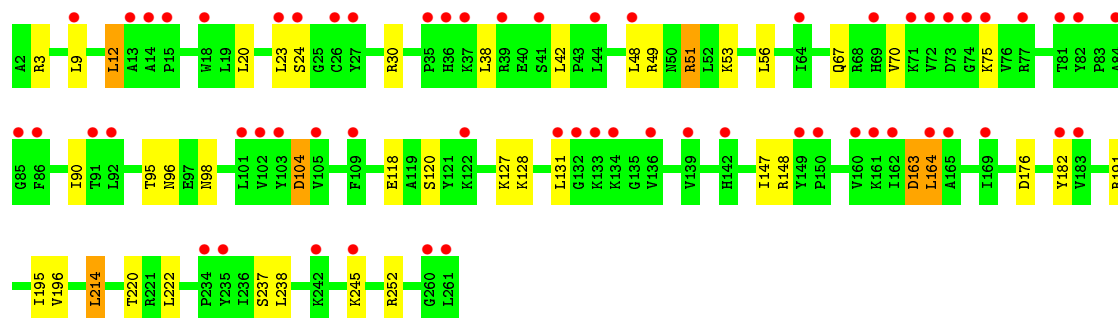
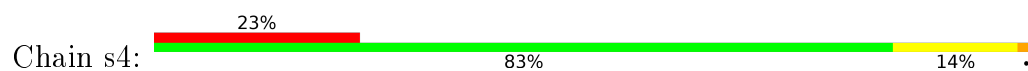




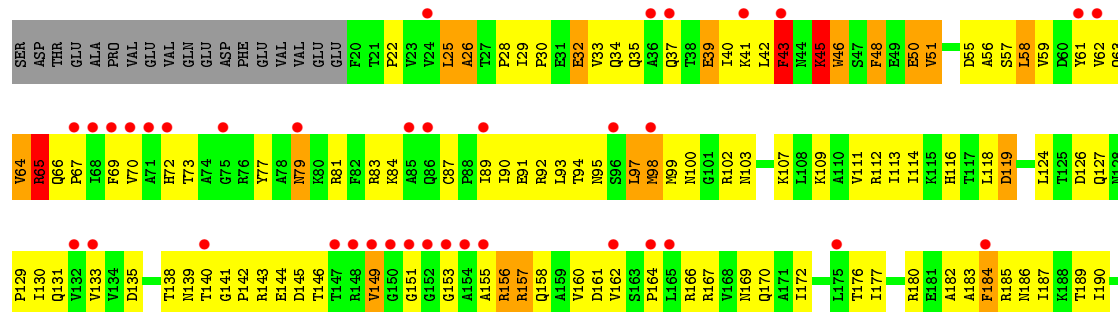
• Molecule 6: 40S ribosomal protein S4-A



• Molecule 6: 40S ribosomal protein S4-A

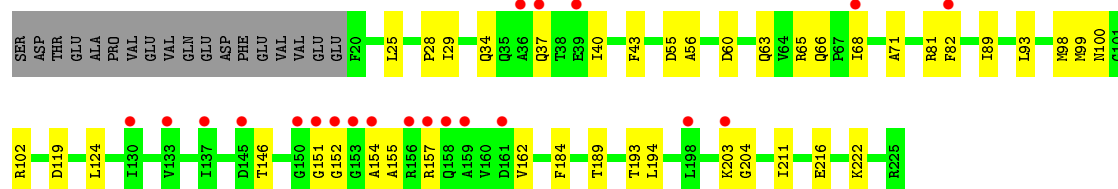
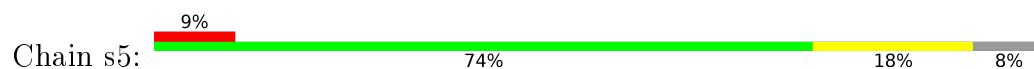


• Molecule 7: 40S ribosomal protein S5

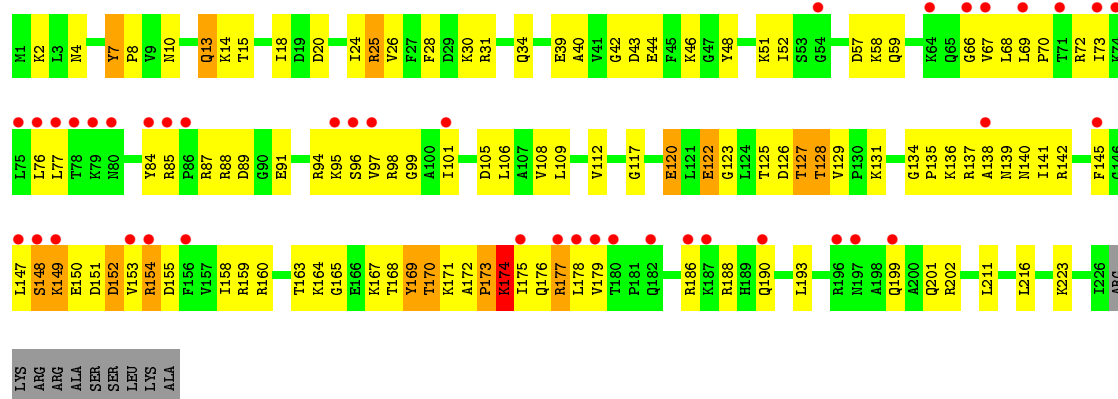




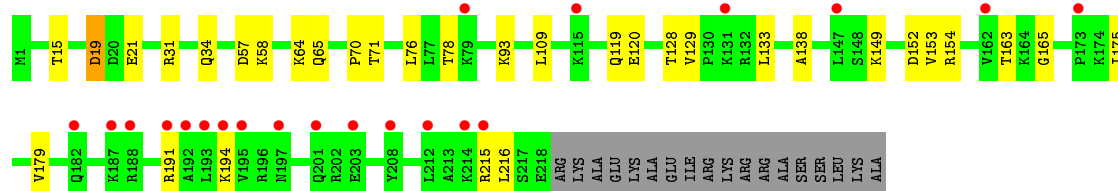
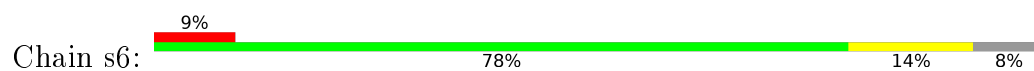
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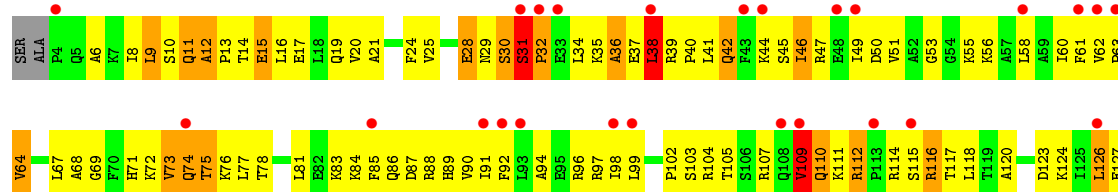
• Molecule 8: 40S ribosomal protein S6-A



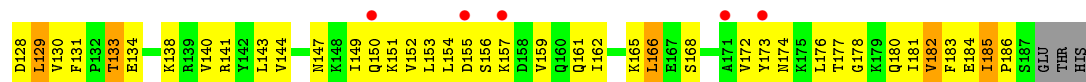
• Molecule 8: 40S ribosomal protein S6-A



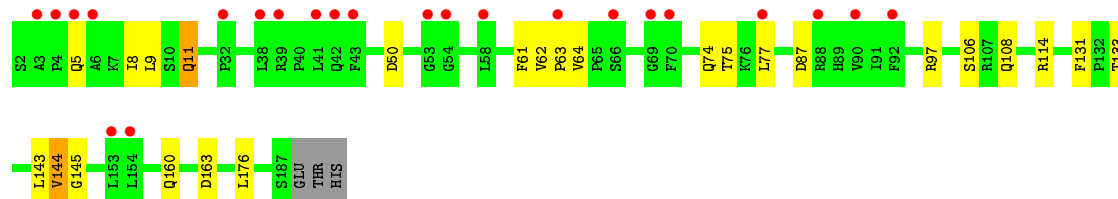
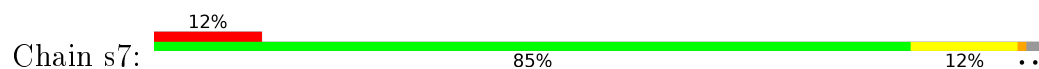
• Molecule 9: 40S ribosomal protein S7-A



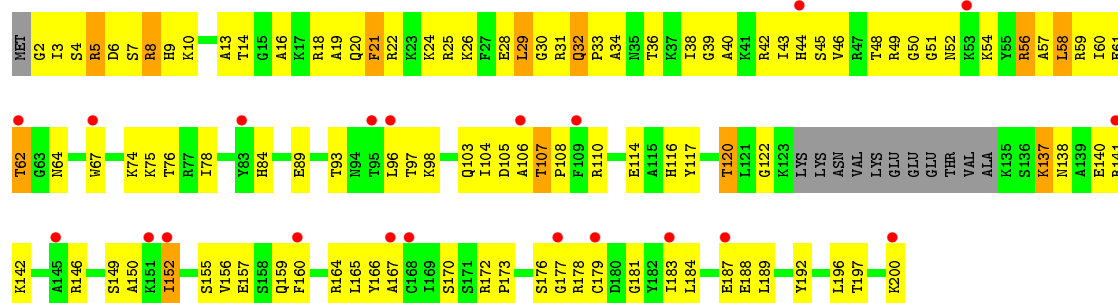
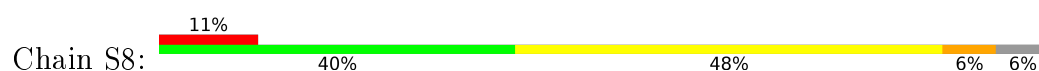




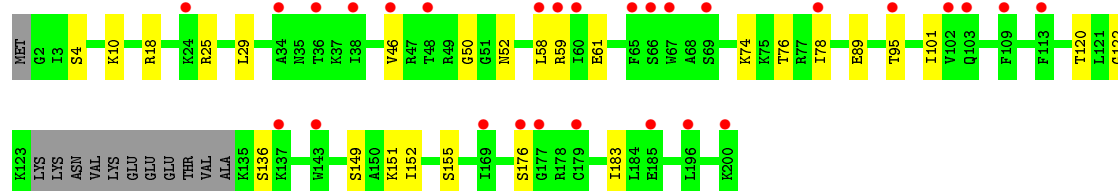
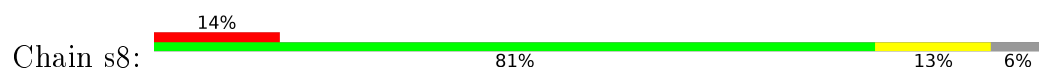
• Molecule 9: 40S ribosomal protein S7-A



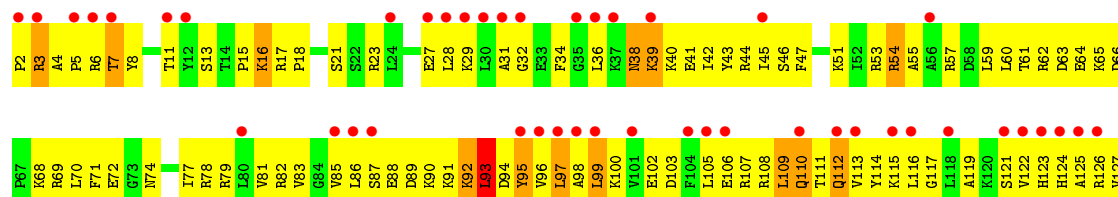
• Molecule 10: 40S ribosomal protein S8-A

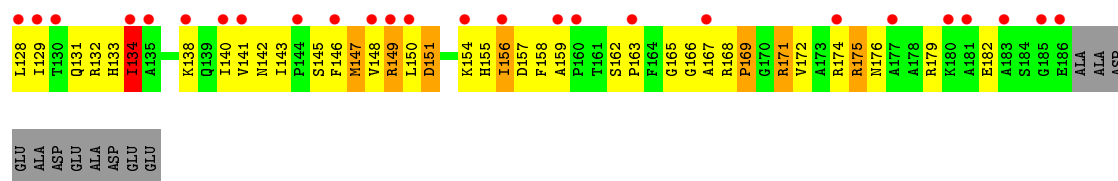


• Molecule 10: 40S ribosomal protein S8-A

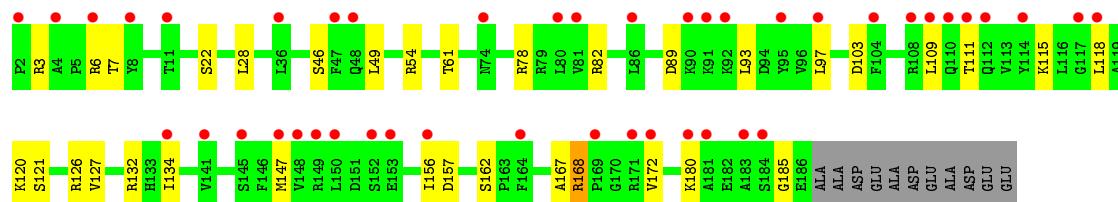
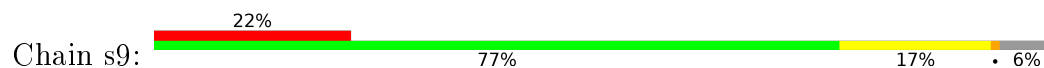


• Molecule 11: 40S ribosomal protein S9-A

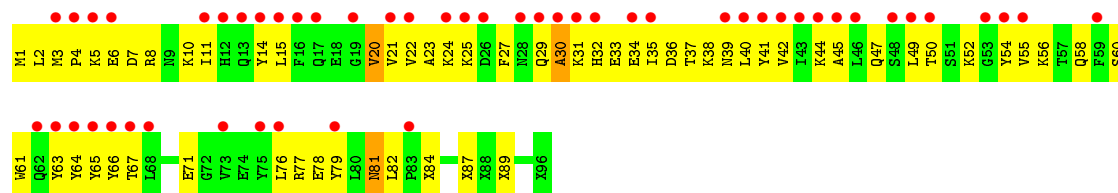




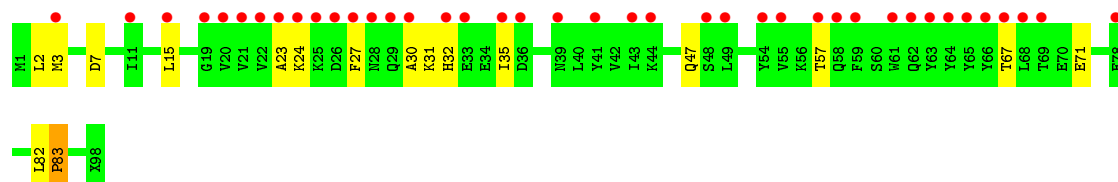
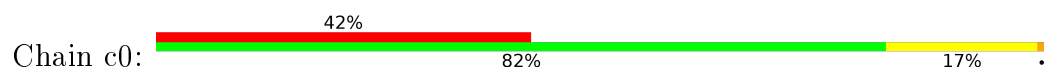
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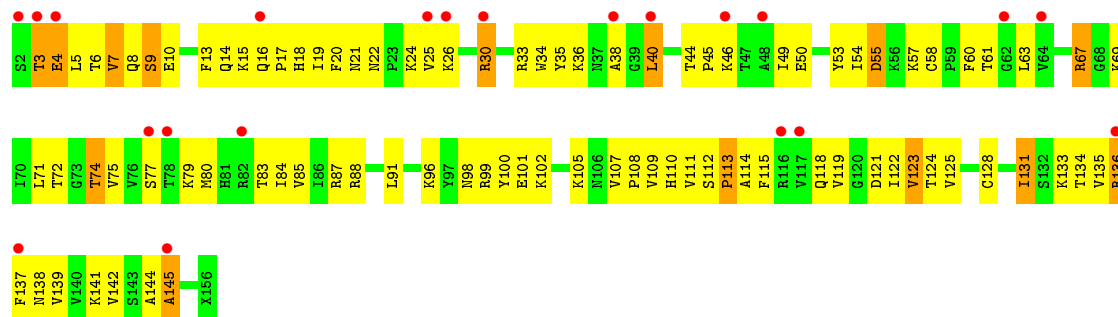
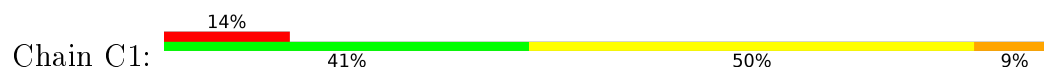
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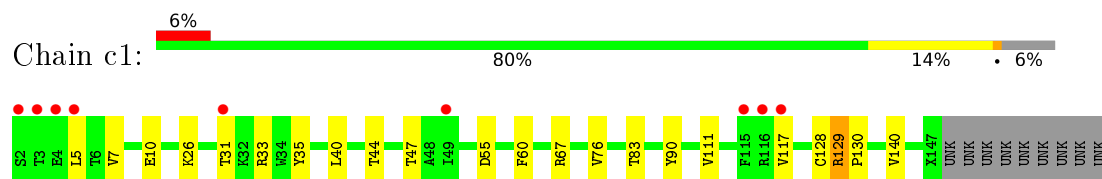
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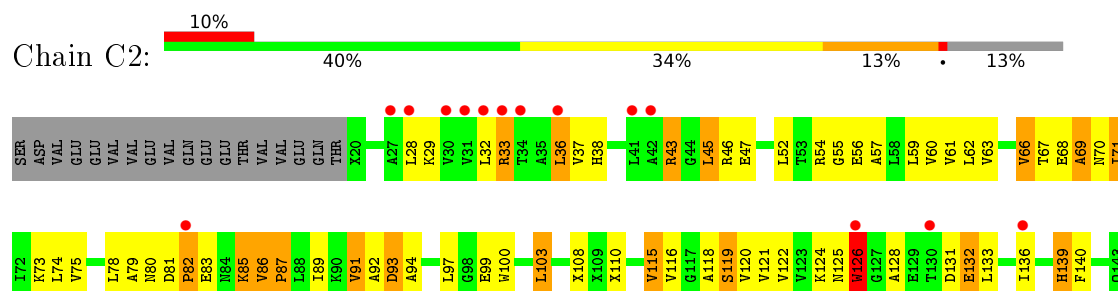
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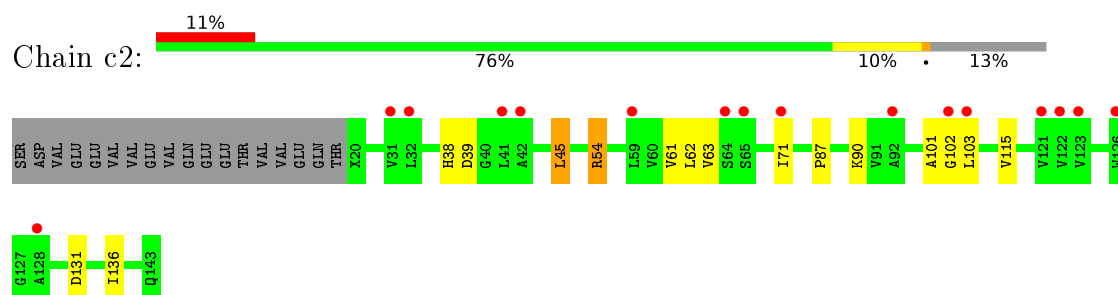
- Molecule 13: 40S ribosomal protein S11-A



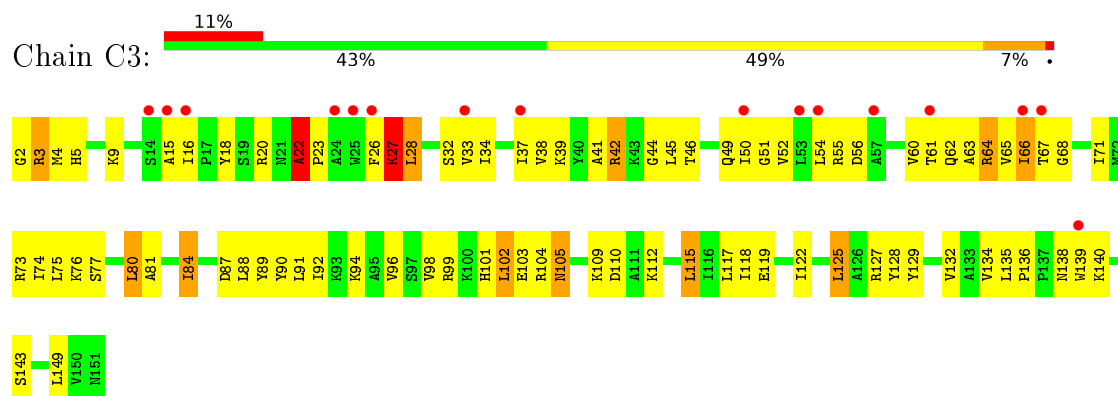
- Molecule 14: 40S ribosomal protein S12



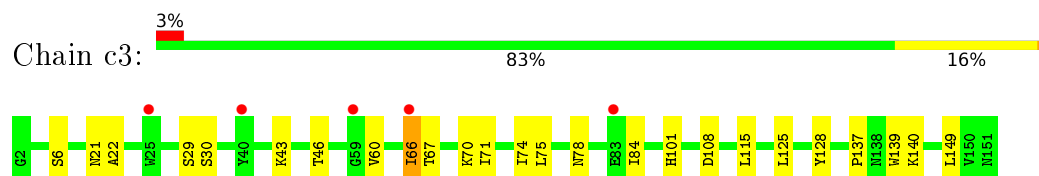
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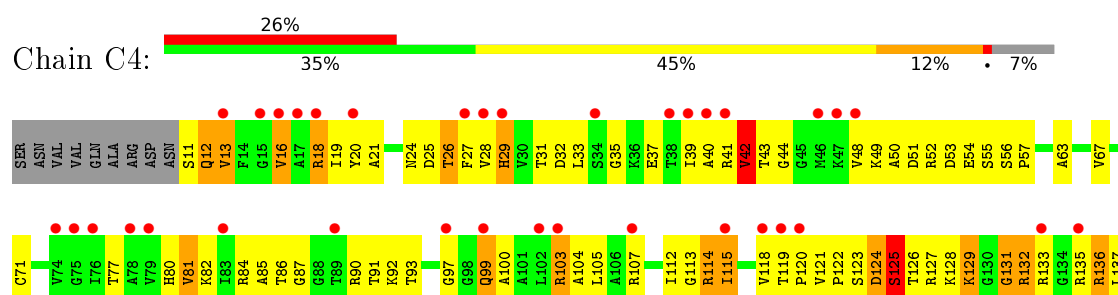
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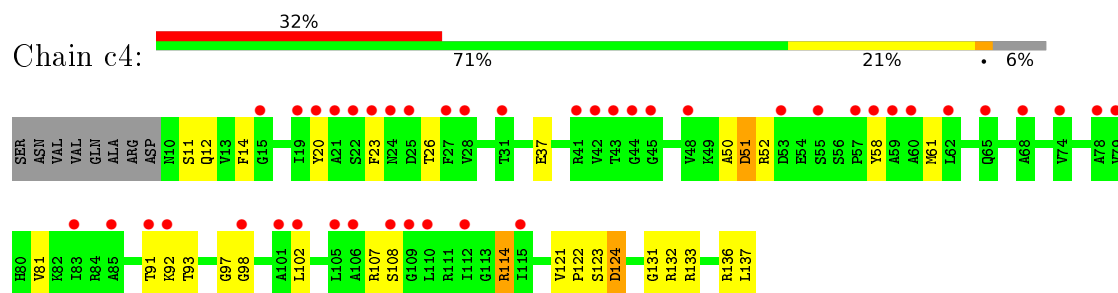
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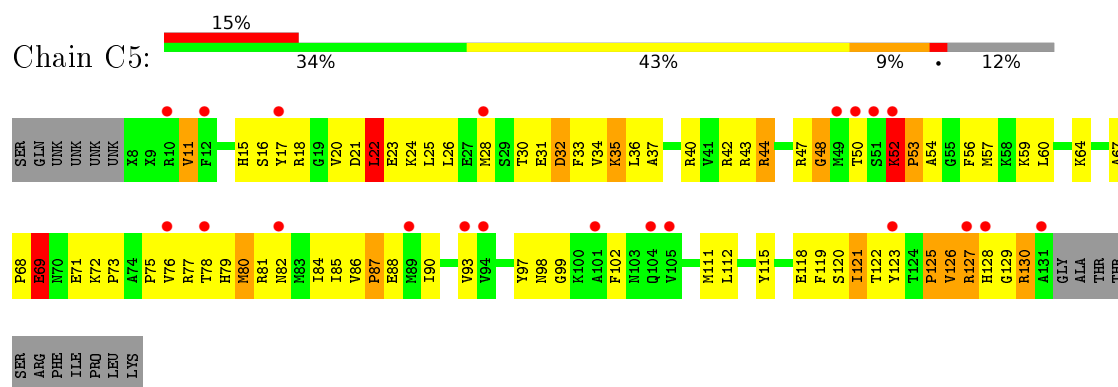
- Molecule 16: 40S ribosomal protein S14-A



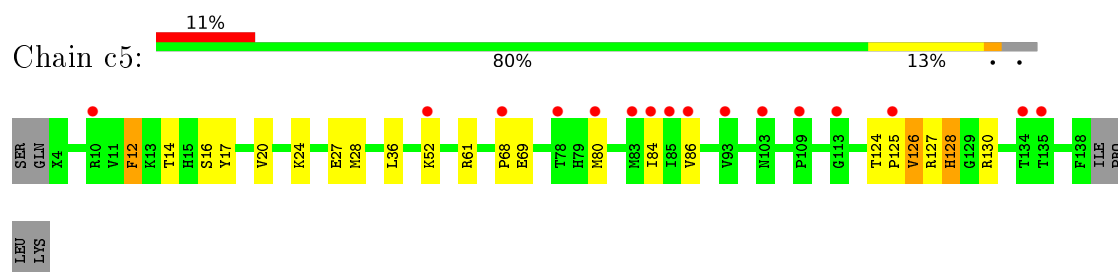
- Molecule 16: 40S ribosomal protein S14-A



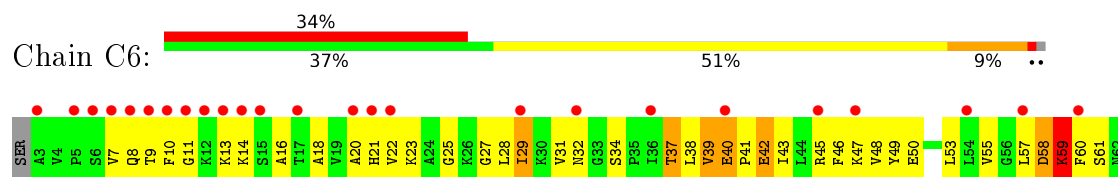
- Molecule 17: 40S ribosomal protein S15 (uS19)

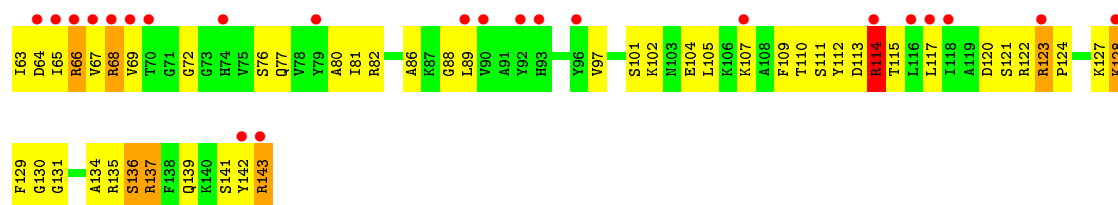


- Molecule 17: 40S ribosomal protein S15 (uS19)

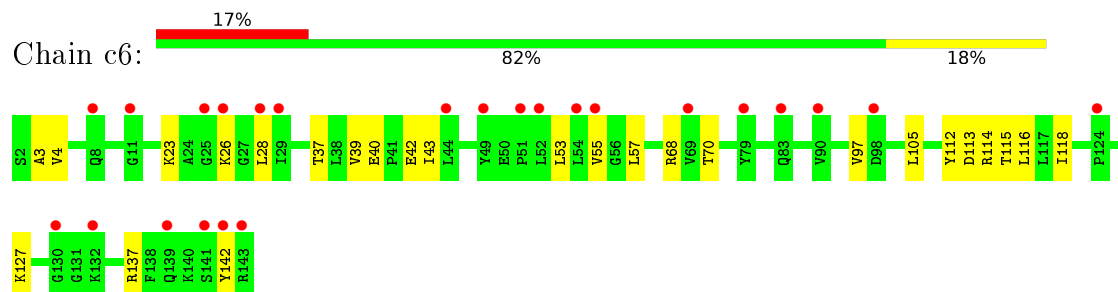


- Molecule 18: 40S ribosomal protein S16-A

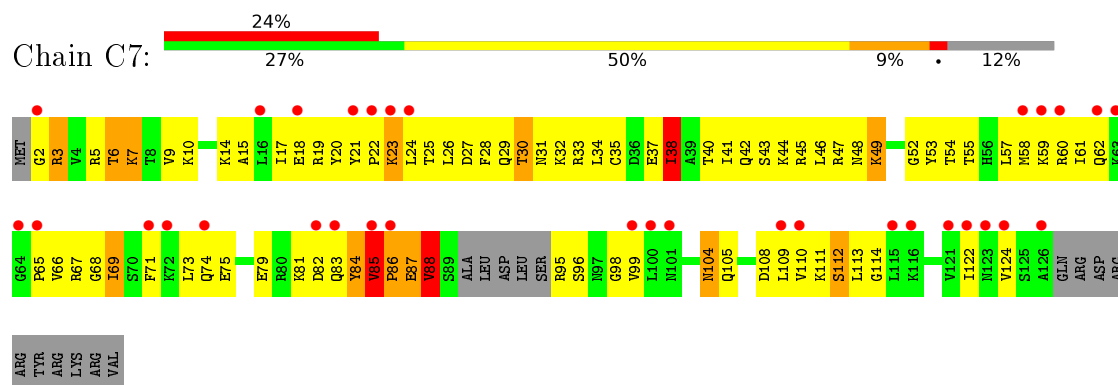




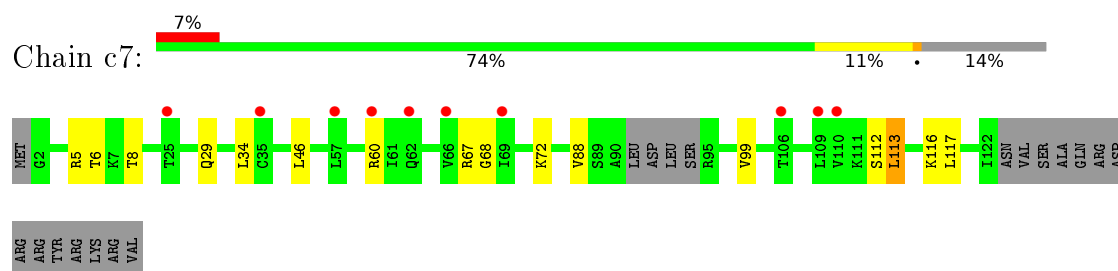
• Molecule 18: 40S ribosomal protein S16-A



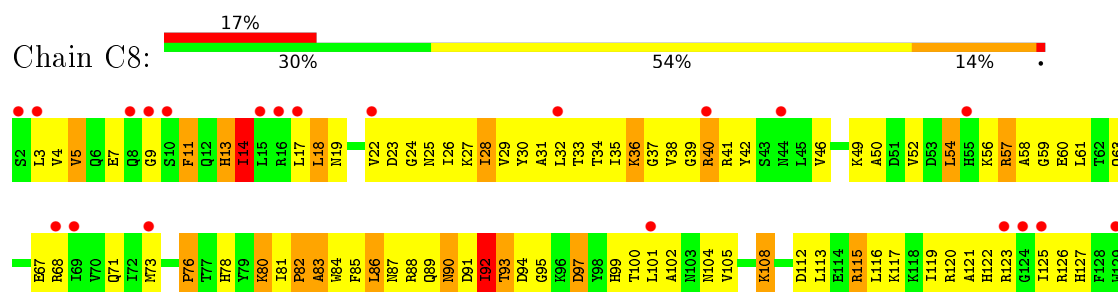
• Molecule 19: 40S ribosomal protein S17-A

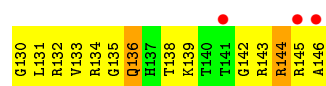


• Molecule 19: 40S ribosomal protein S17-A

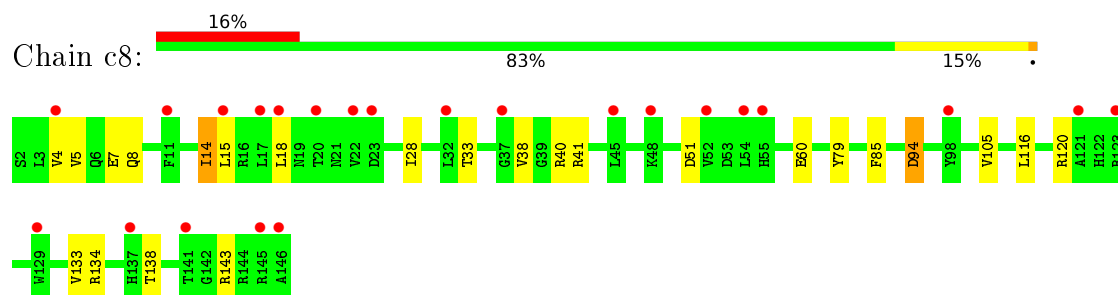


• Molecule 20: 40S ribosomal protein S18-A

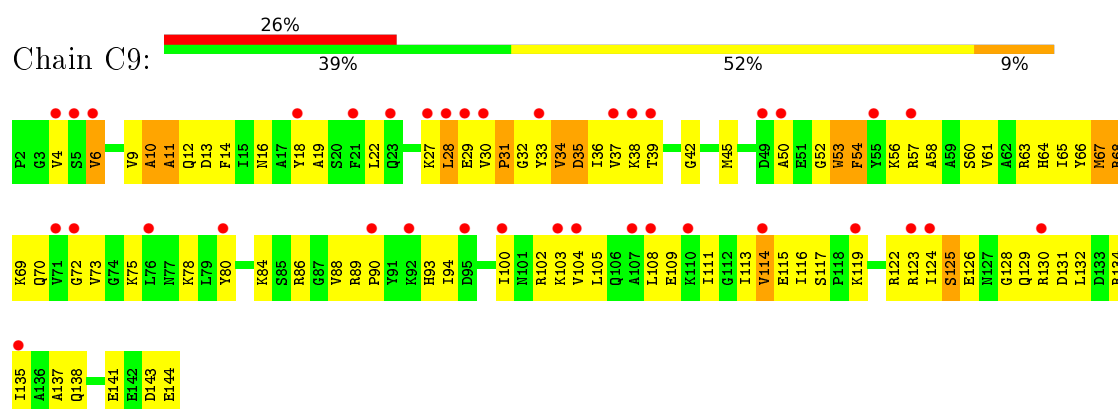




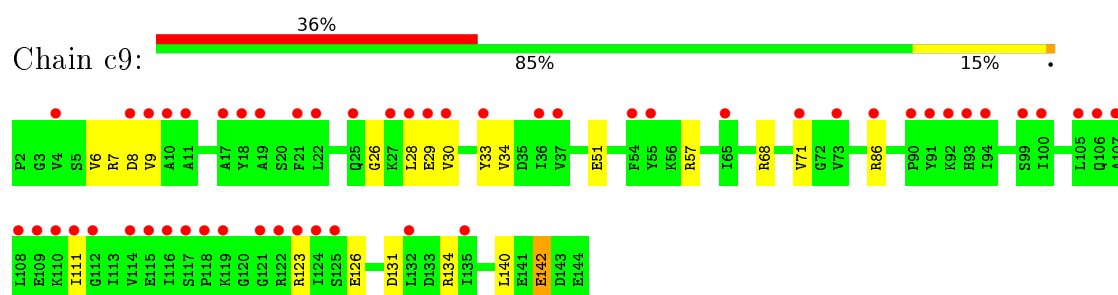
- Molecule 20: 40S ribosomal protein S18-A



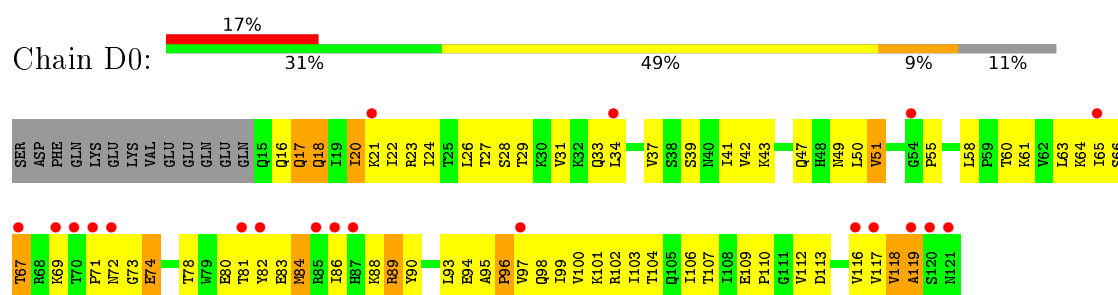
- Molecule 21: 40S ribosomal protein S19-A



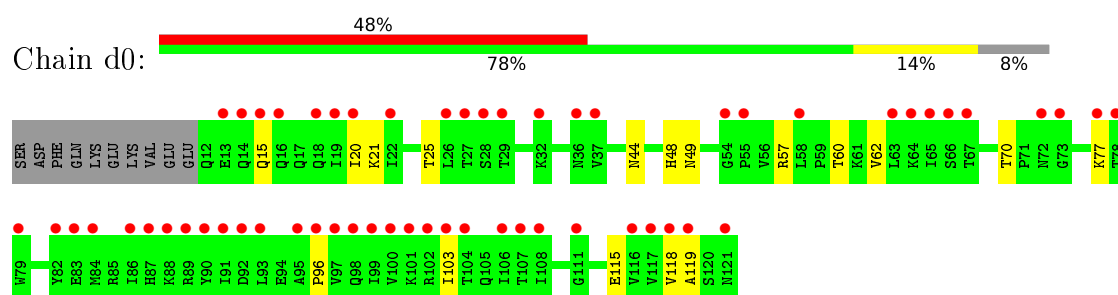
- Molecule 21: 40S ribosomal protein S19-A



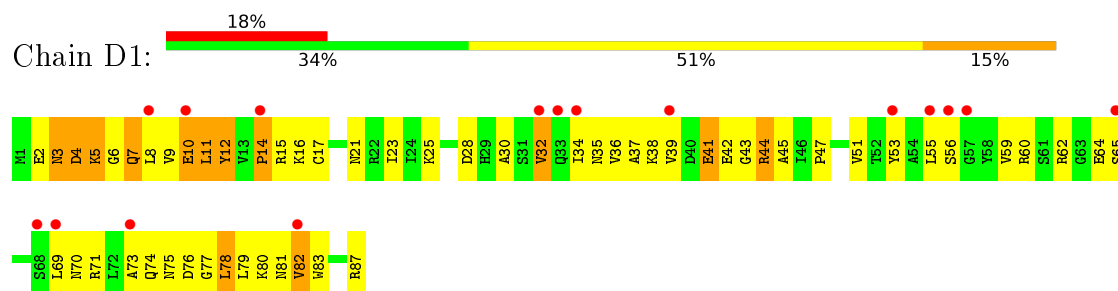
- Molecule 22: 40S ribosomal protein S20



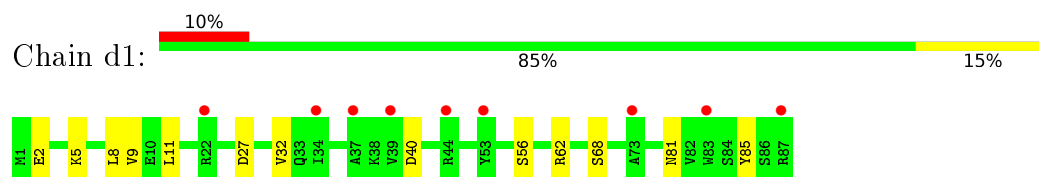
- Molecule 22: 40S ribosomal protein S20



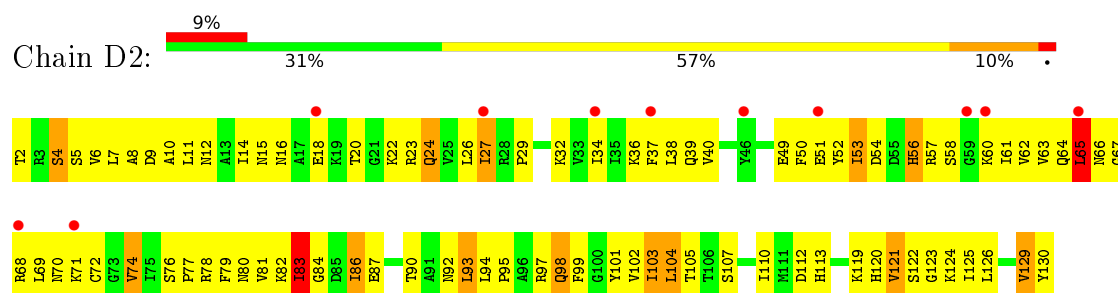
- Molecule 23: 40S ribosomal protein S21-A



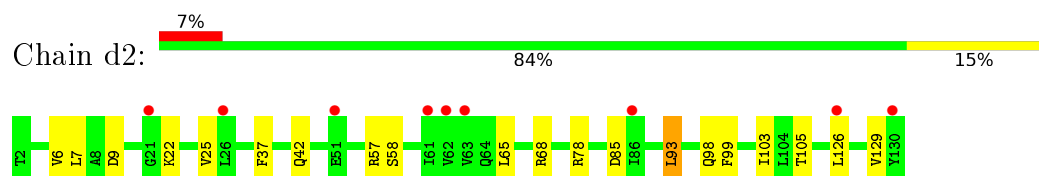
- Molecule 23: 40S ribosomal protein S21-A



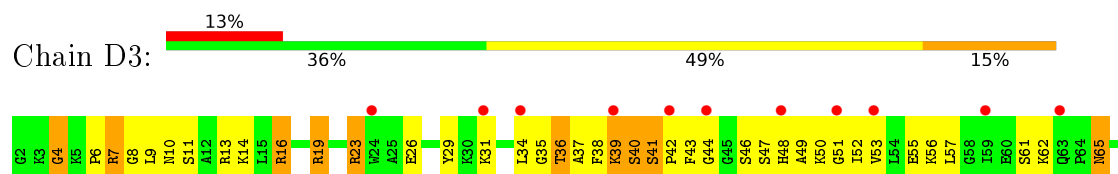
- Molecule 24: 40S ribosomal protein S22-A

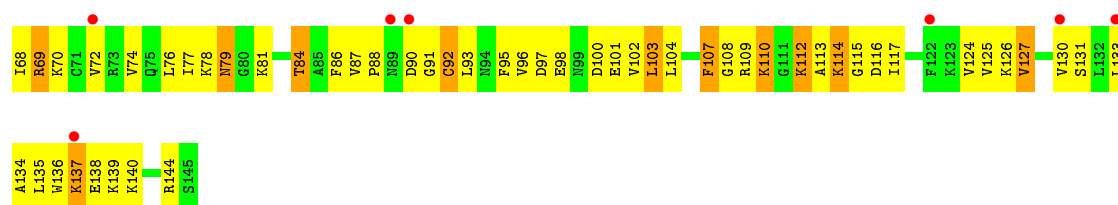


- Molecule 24: 40S ribosomal protein S22-A

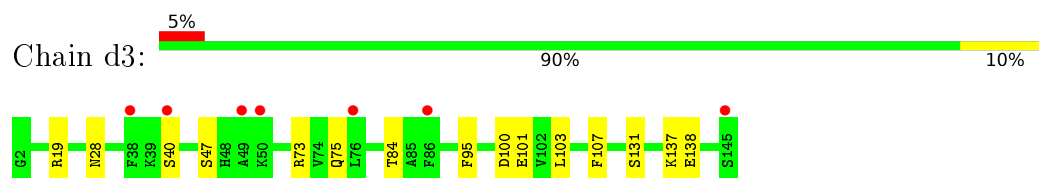


- Molecule 25: 40S ribosomal protein S23-A

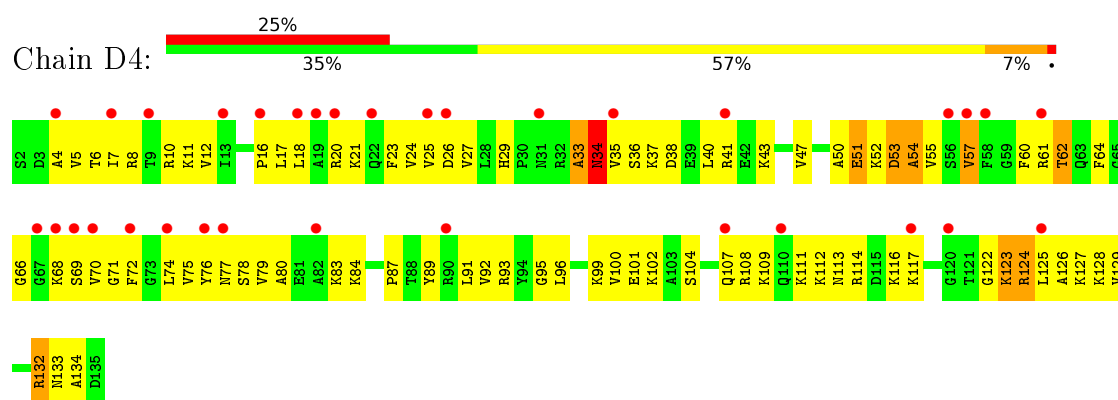




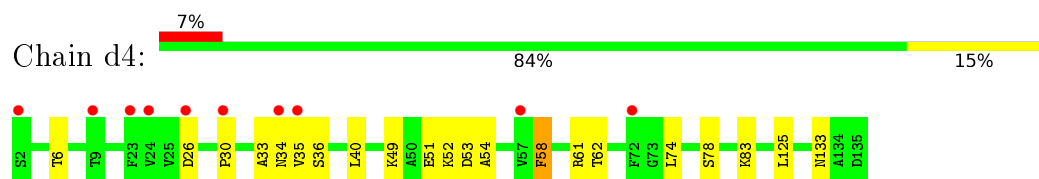
- Molecule 25: 40S ribosomal protein S23-A



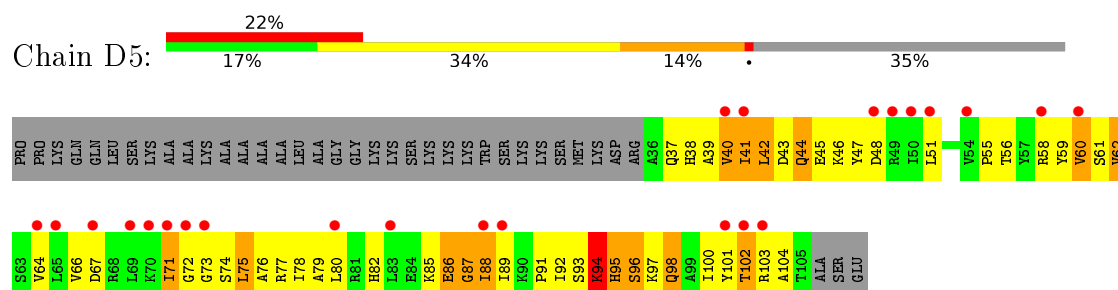
- Molecule 26: 40S ribosomal protein S24-A



- Molecule 26: 40S ribosomal protein S24-A



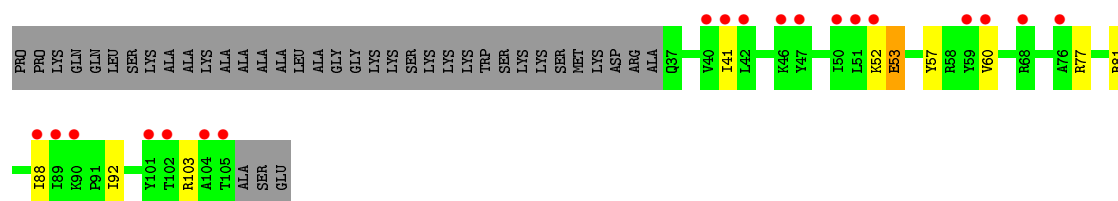
- Molecule 27: 40S ribosomal protein S25-A



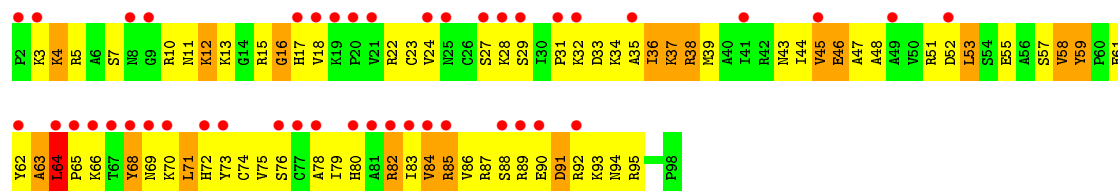
- Molecule 27: 40S ribosomal protein S25-A



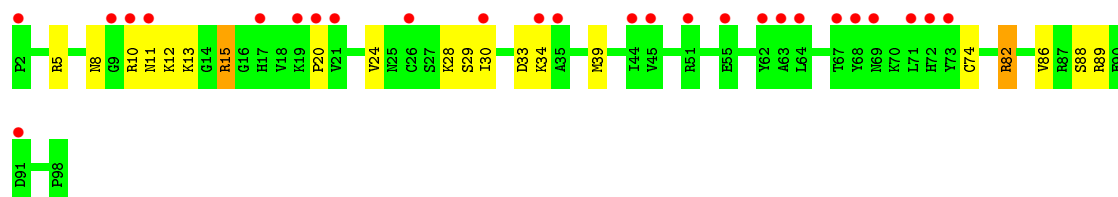
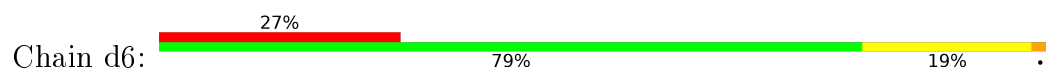




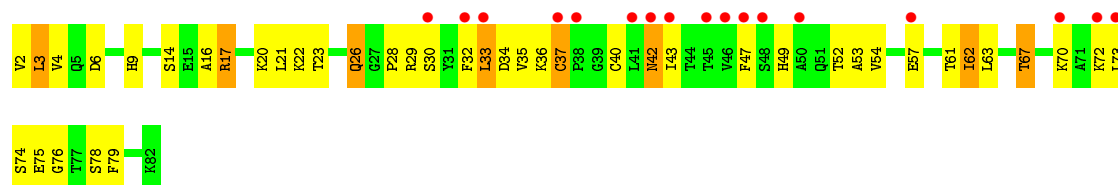
• Molecule 28: 40S ribosomal protein S26-B



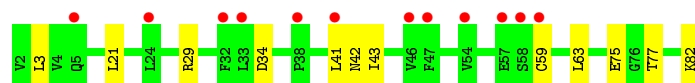
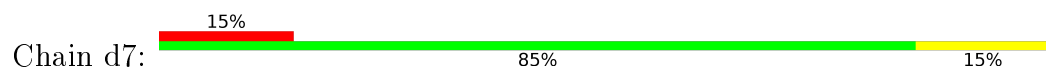
• Molecule 28: 40S ribosomal protein S26-B



• Molecule 29: 40S ribosomal protein S27-A

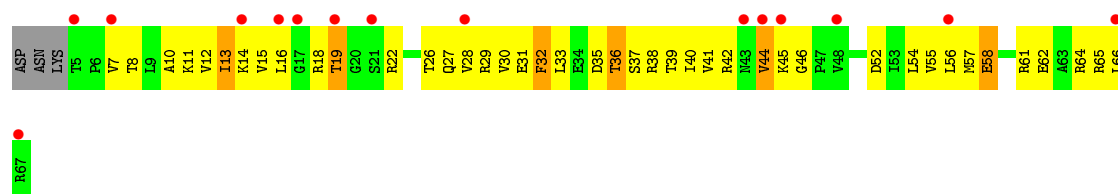


• Molecule 29: 40S ribosomal protein S27-A

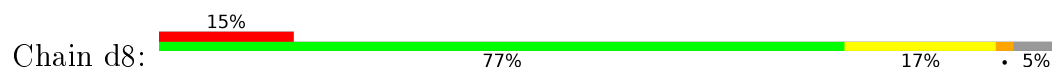


• Molecule 30: 40S ribosomal protein S28-A





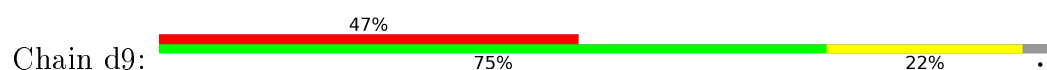
- Molecule 30: 40S ribosomal protein S28-A



- Molecule 31: 40S ribosomal protein S29-A



- Molecule 31: 40S ribosomal protein S29-A



- Molecule 32: 40S ribosomal protein S30-A

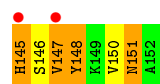


- Molecule 32: 40S ribosomal protein S30-A

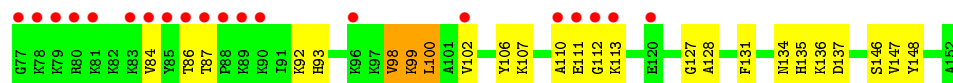


- Molecule 33: Ubiquitin-40S ribosomal protein S31

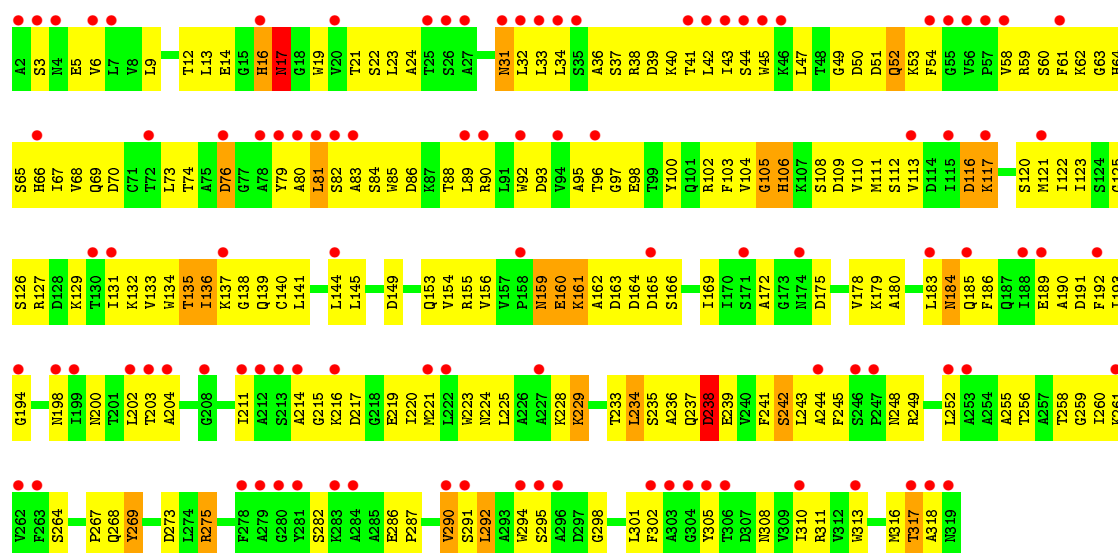




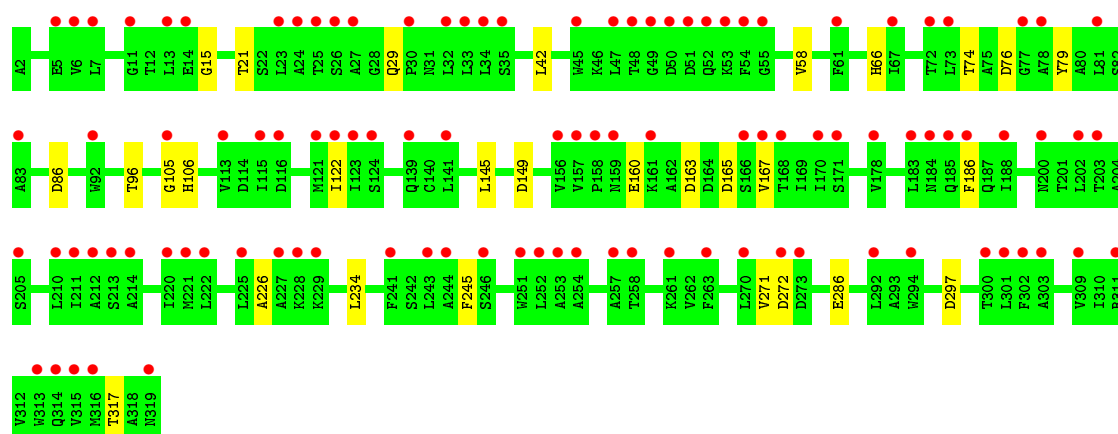
- Molecule 33: Ubiquitin-40S ribosomal protein S31



- Molecule 34: Guanine nucleotide-binding protein subunit beta-like protein



- Molecule 34: Guanine nucleotide-binding protein subunit beta-like protein

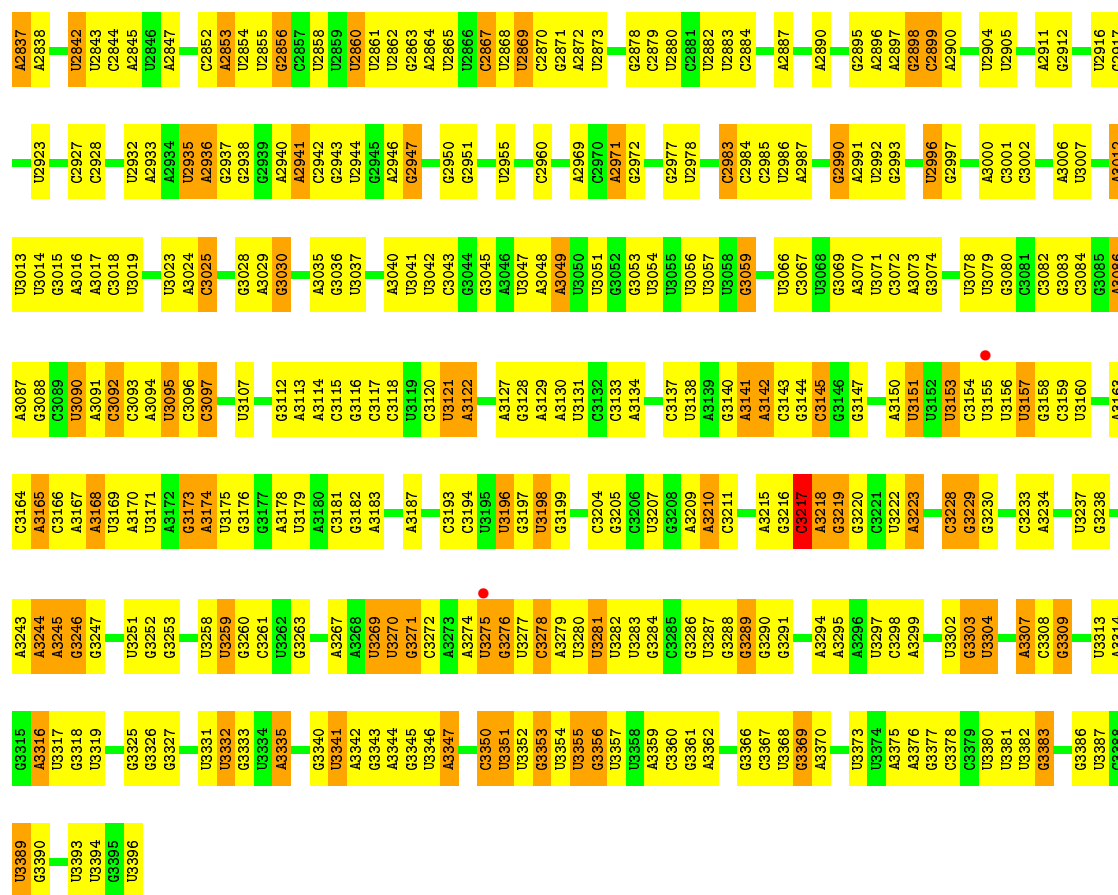


- Molecule 35: Suppressor protein STM1

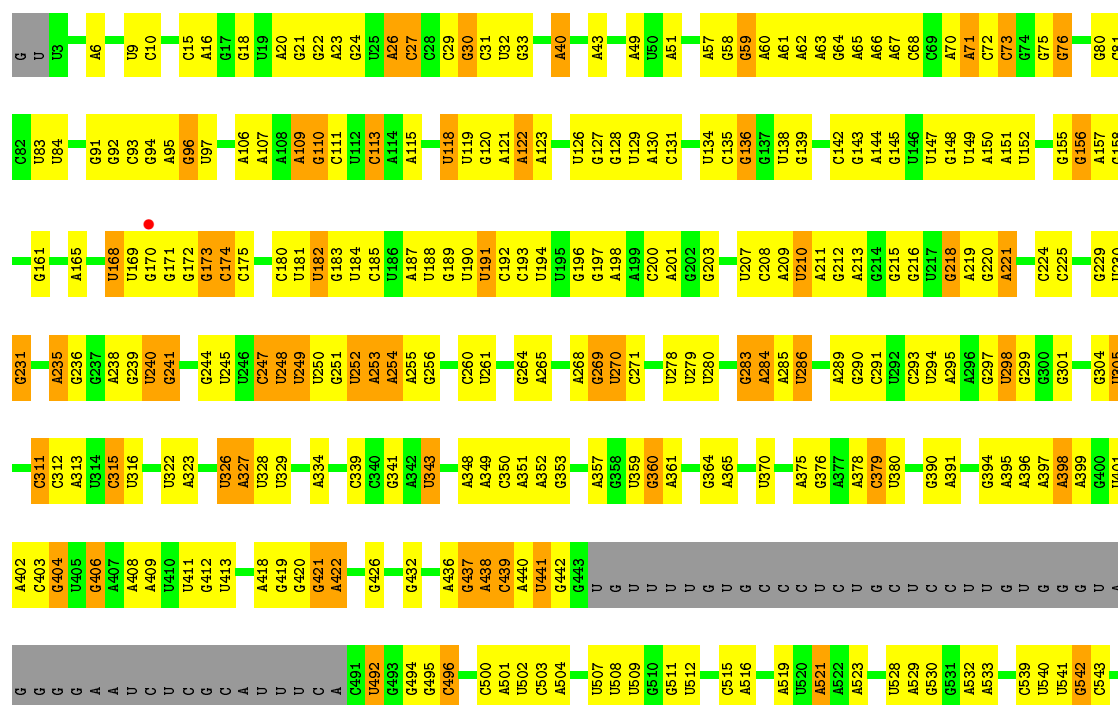
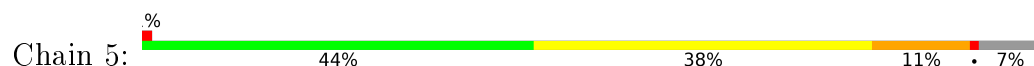








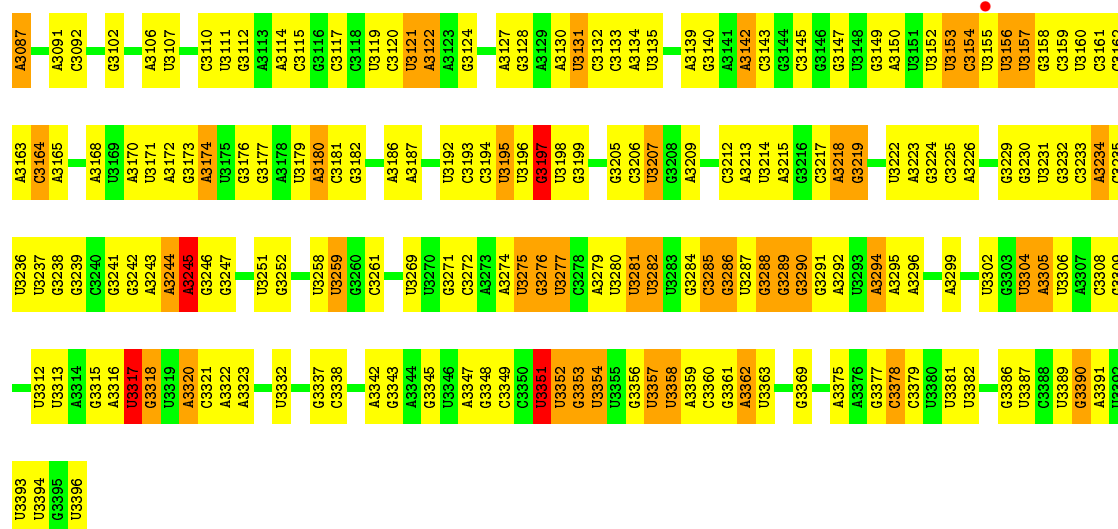
• Molecule 36: 25S ribosomal RNA



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G1829	G1747	C1657	C1578	G1488	A1399	A1129	C1049	A973	A884	U777	C695	A621	G547
G1830		G1658	A1580	A1489	G1400	A1130	C1048	G974	C986		C696	A622	G548
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A1842	C1762			U1500	G1421	G1145	U1070		A906	U790	A706	G634	U558
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A1787	G1713		U1544		A1449	A1263	U1097	G1012		G826	A735	G659	A589
	A1714		G1544		G1450	A1353	A1098	G1013	U942	A830	A736	G661	G590
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A1879	G1628				A1461	A1271	G1101	C1016		G838	C743	U664	C593
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			A1637		U1471	G1285	A1112	A1027	C957	U855	A761	G678	A603
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U1916	U1740		C1574		U1483	A1391	G1213	U1041		G881		A693	
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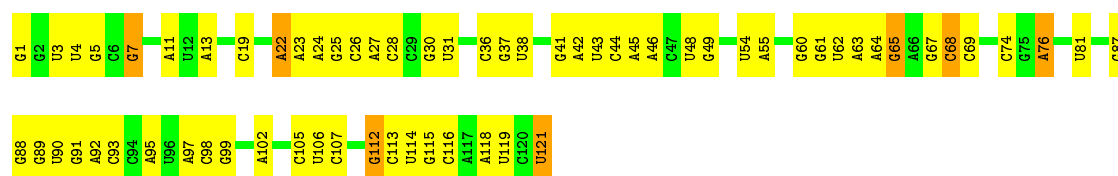


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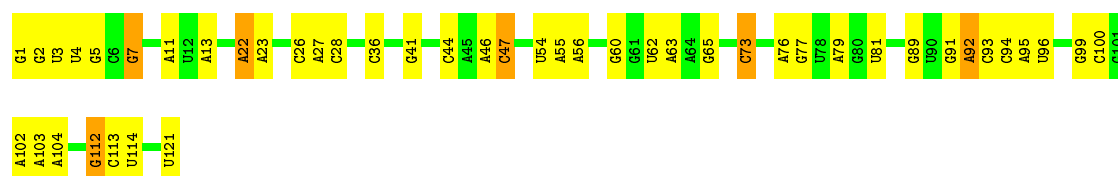
• Molecule 37: 5S ribosomal RNA

Chain 3: 46% 48% 6%



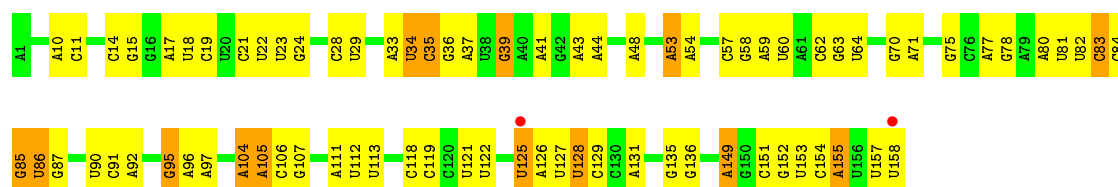
• Molecule 37: 5S ribosomal RNA

Chain 7: 62% 33% 5%



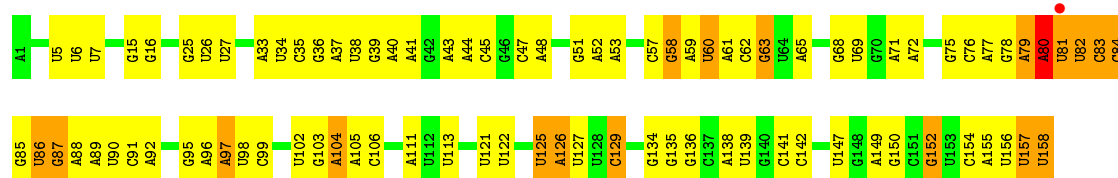
• Molecule 38: 5.8S ribosomal RNA

Chain 4: 51% 41% 9%

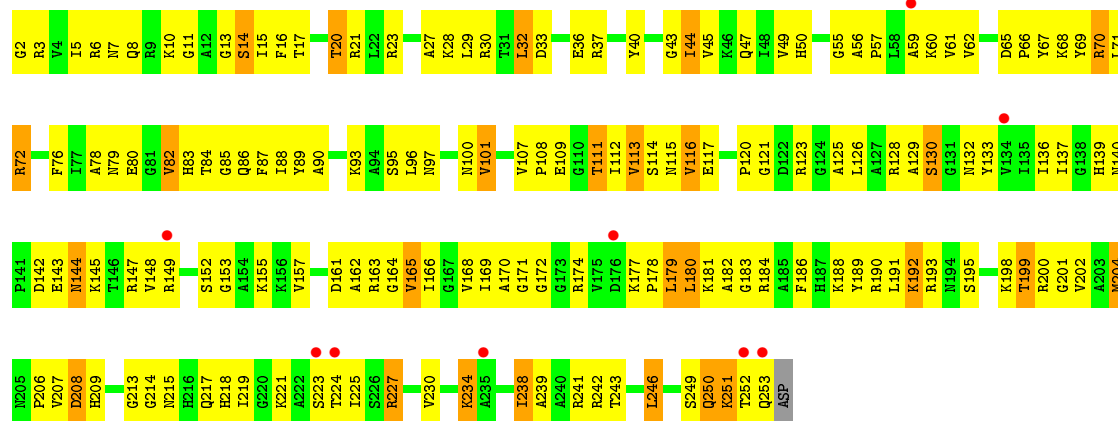


• Molecule 38: 5.8S ribosomal RNA

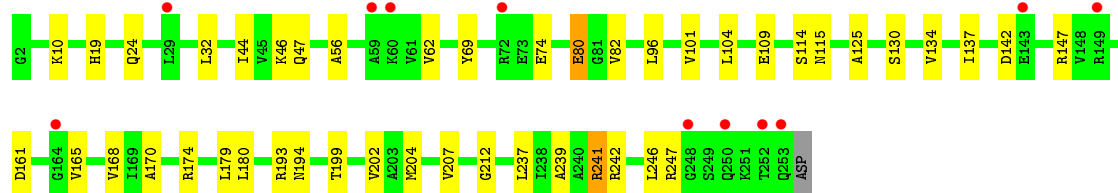
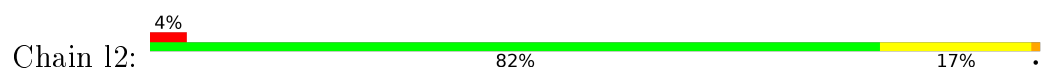
Chain 8: 44% 44% 11%



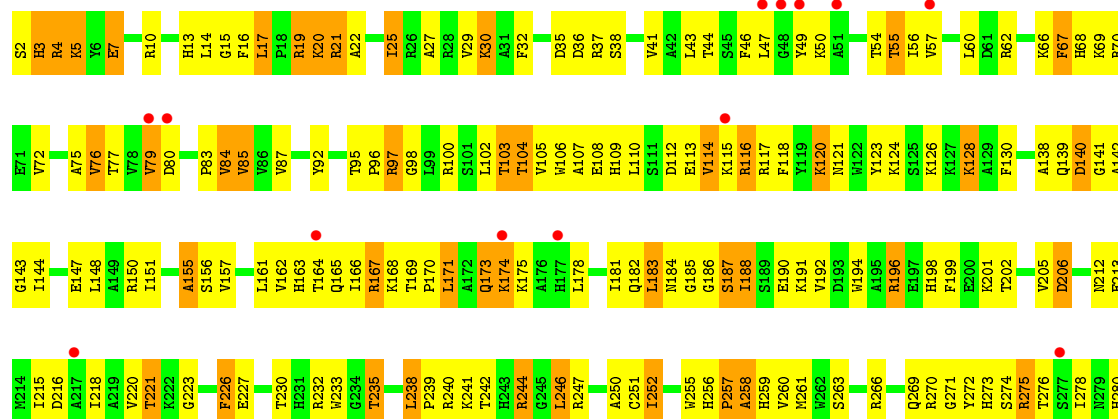
• Molecule 39: 60S ribosomal protein L2-A

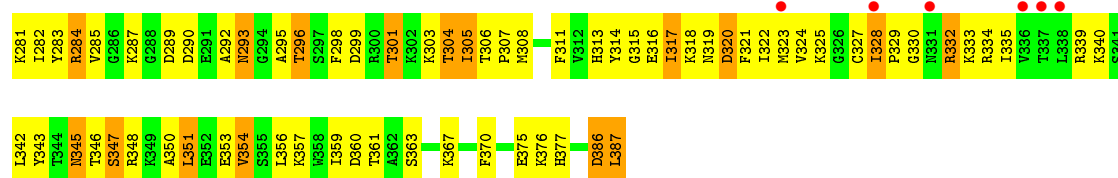


• Molecule 39: 60S ribosomal protein L2-A

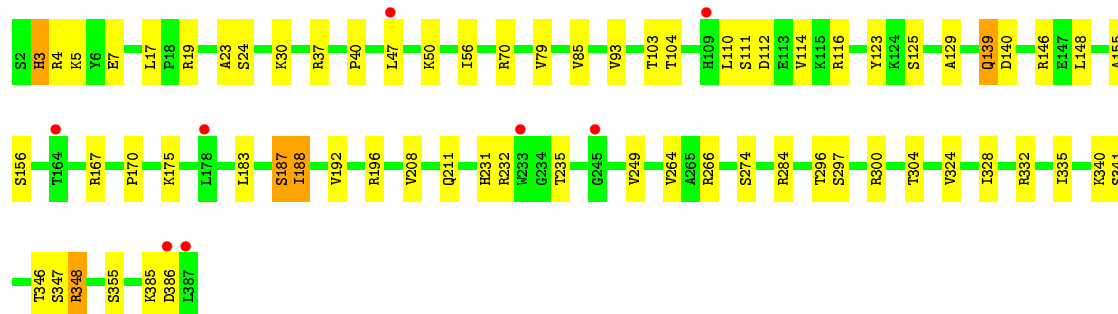
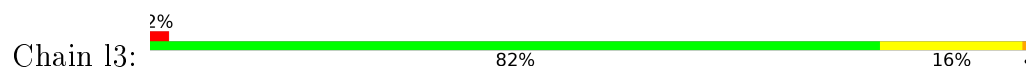


• Molecule 40: 60S ribosomal protein L3

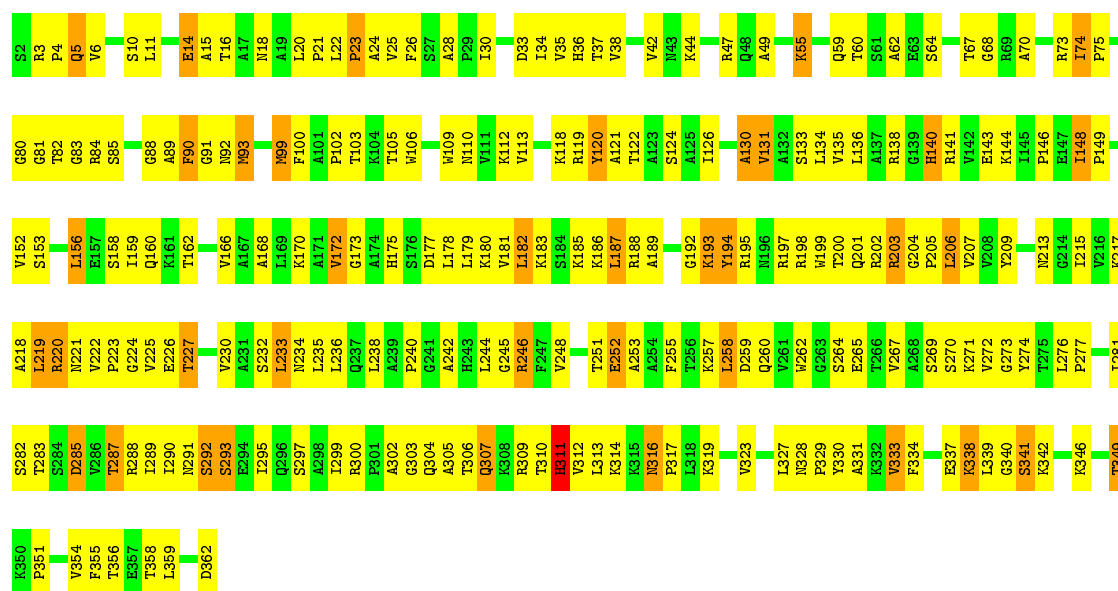




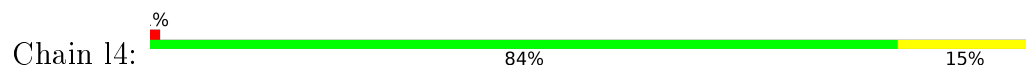
• Molecule 40: 60S ribosomal protein L3

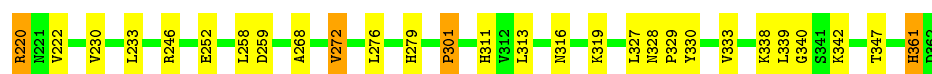


• Molecule 41: 60S ribosomal protein L4-A

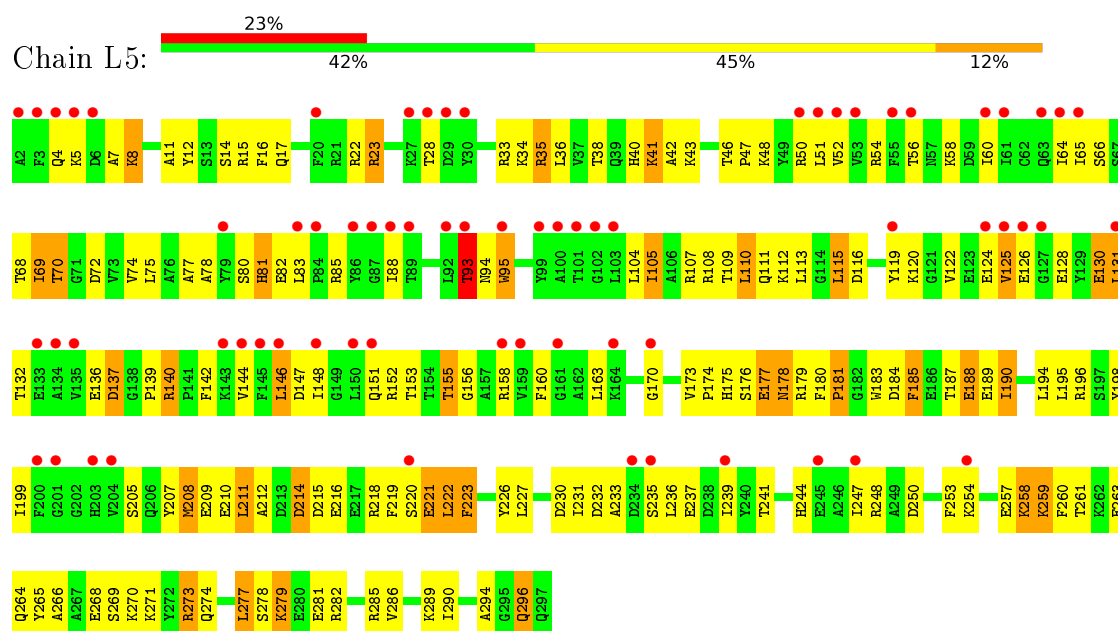


• Molecule 41: 60S ribosomal protein L4-A

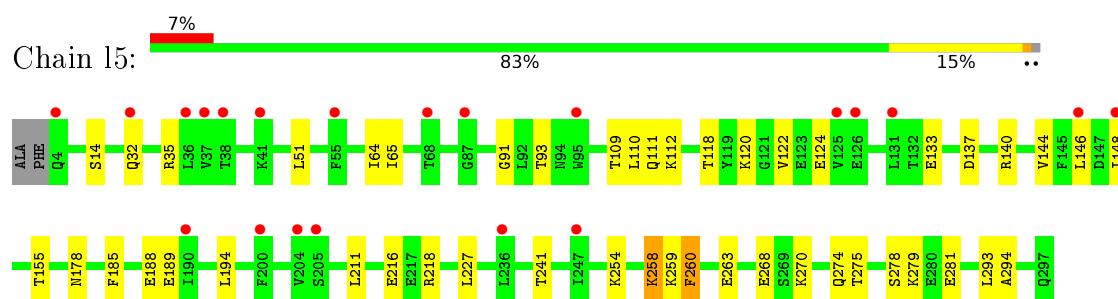




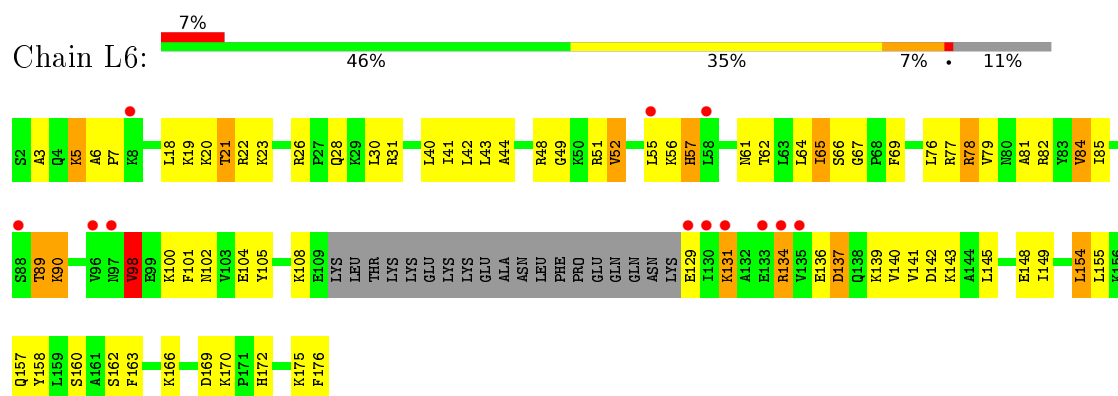
• Molecule 42: 60S ribosomal protein L5



• Molecule 42: 60S ribosomal protein L5

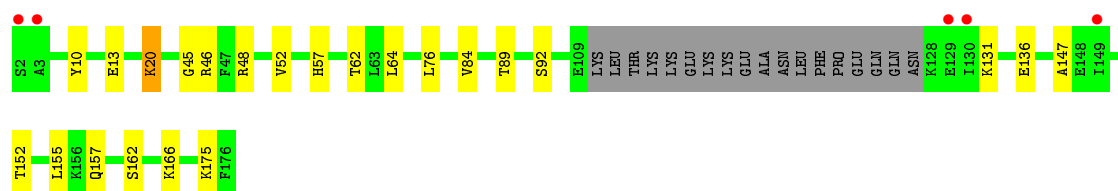


• Molecule 43: 60S ribosomal protein L6-A

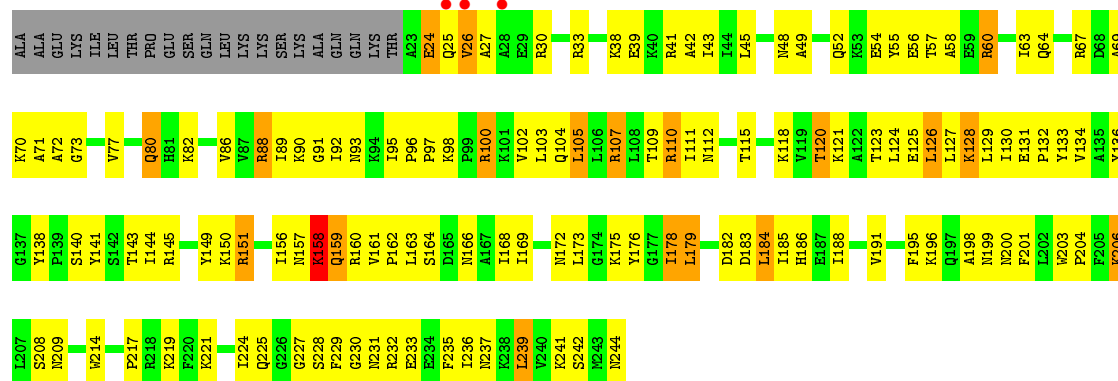


• Molecule 43: 60S ribosomal protein L6-A

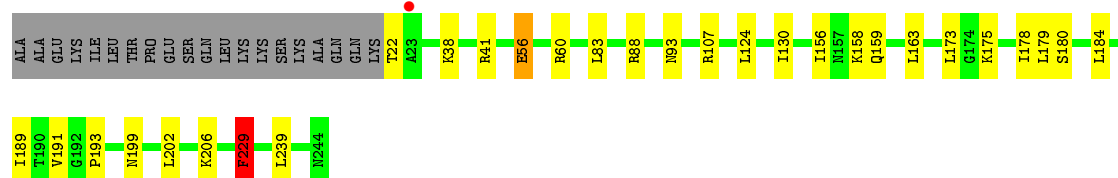
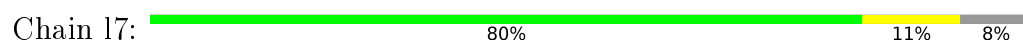




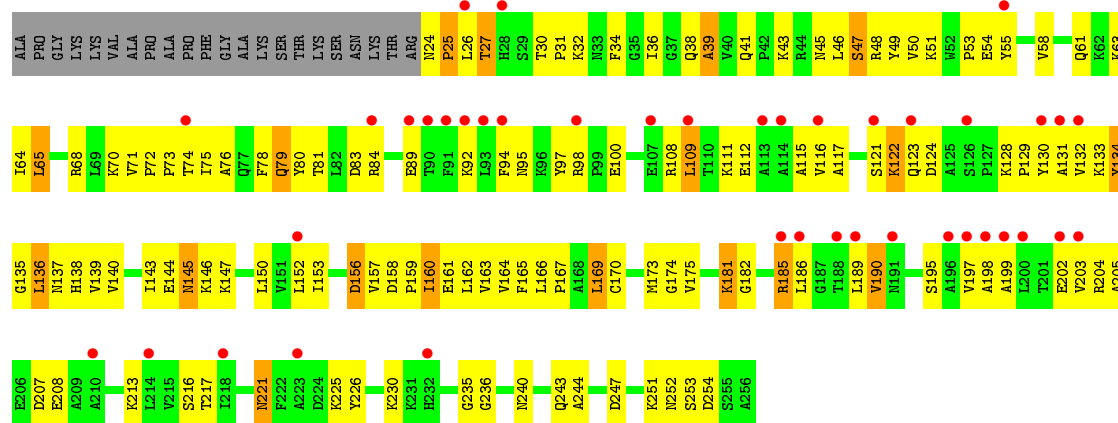
• Molecule 44: 60S ribosomal protein L7-A



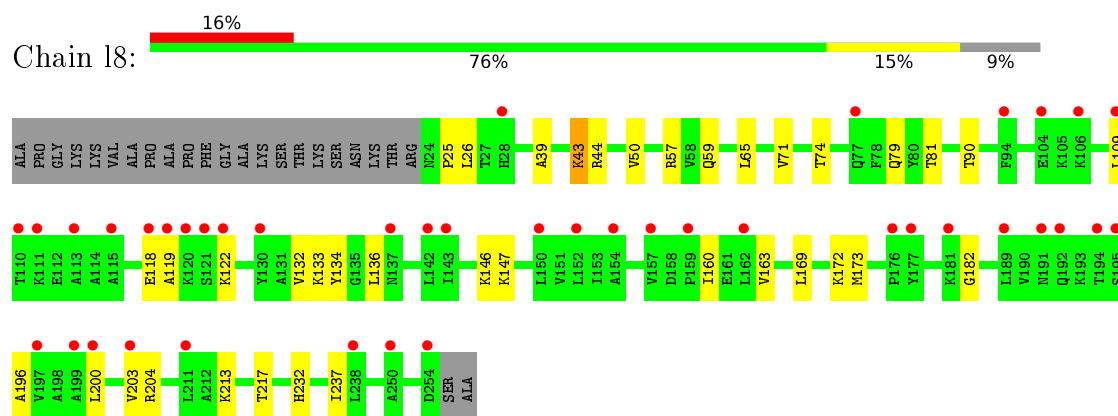
• Molecule 44: 60S ribosomal protein L7-A



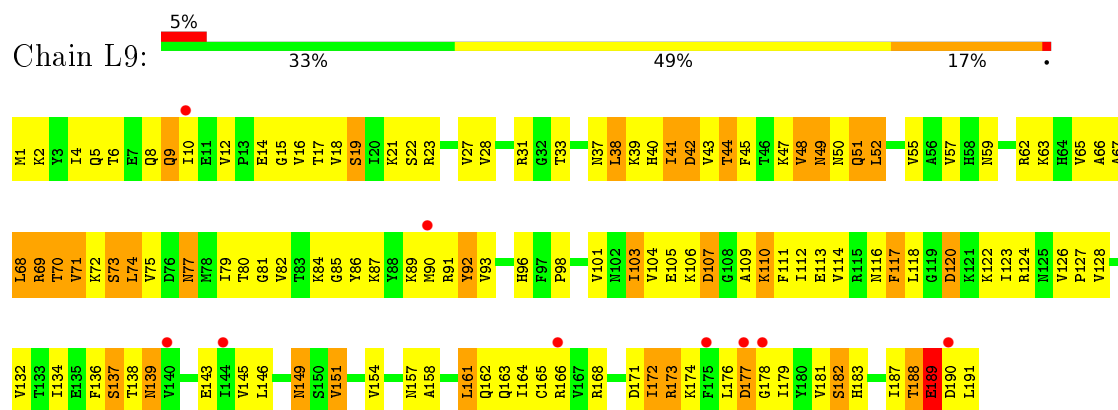
• Molecule 45: 60S ribosomal protein L8-A



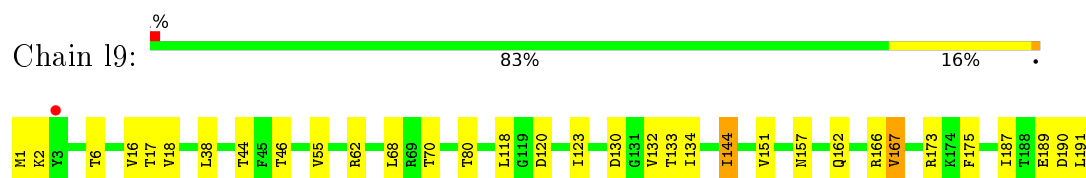
• Molecule 45: 60S ribosomal protein L8-A



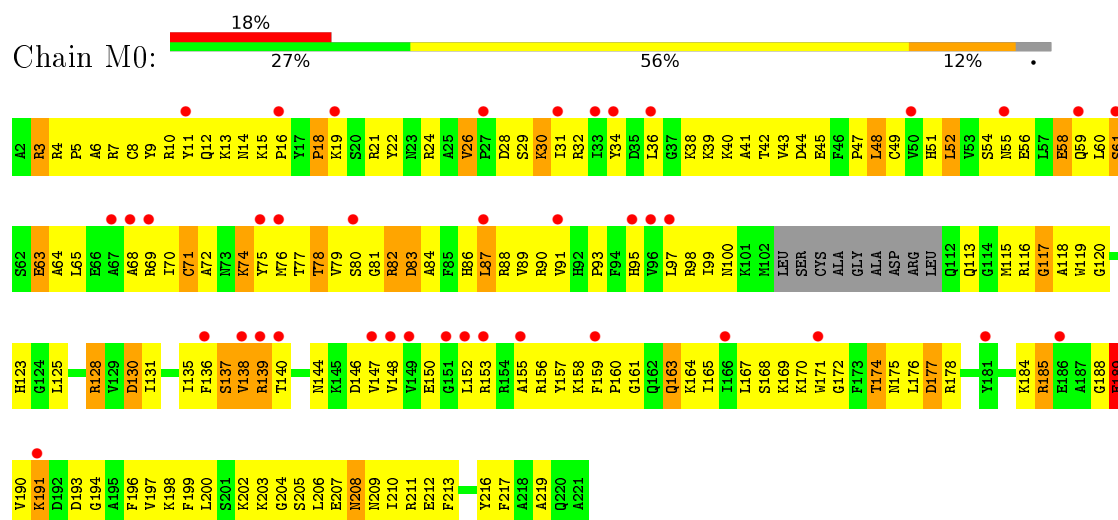
- Molecule 46: 60S ribosomal protein L9-A



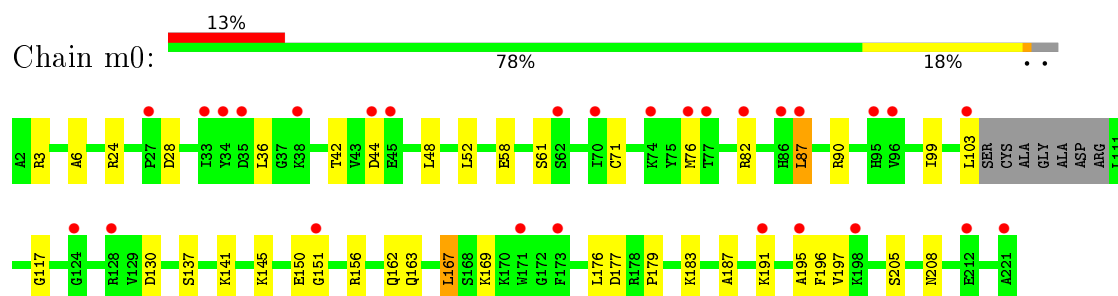
- Molecule 46: 60S ribosomal protein L9-A



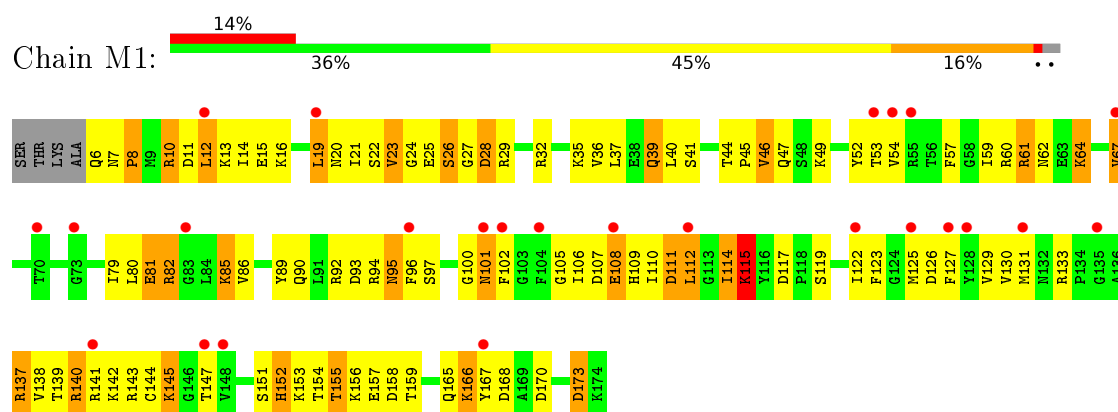
- Molecule 47: 60S ribosomal protein L10



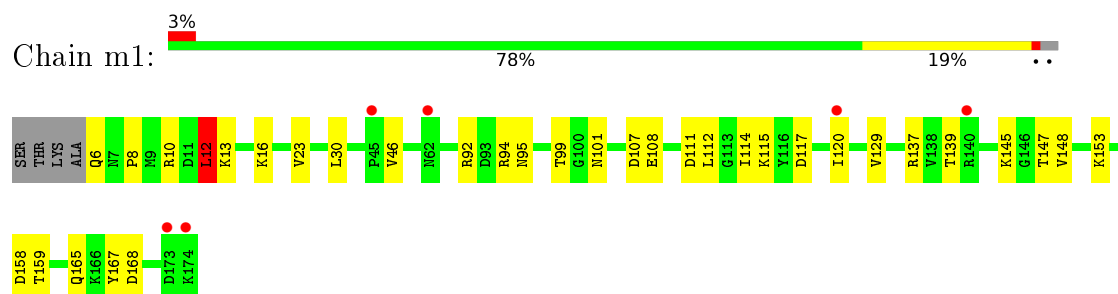
- Molecule 47: 60S ribosomal protein L10



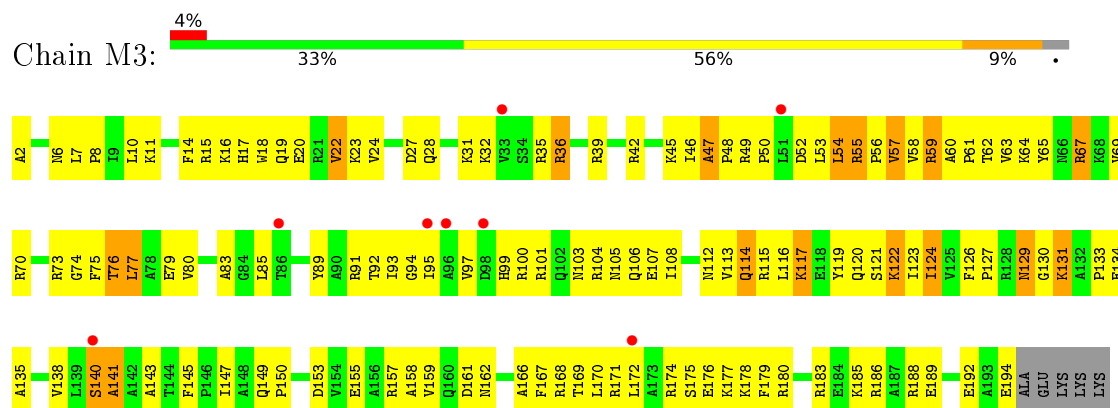
- Molecule 48: 60S ribosomal protein L11-B



- Molecule 48: 60S ribosomal protein L11-B

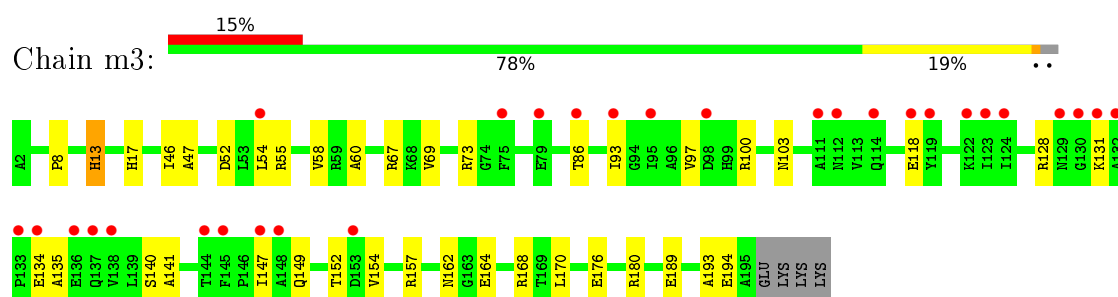


- Molecule 49: 60S ribosomal protein L13-A

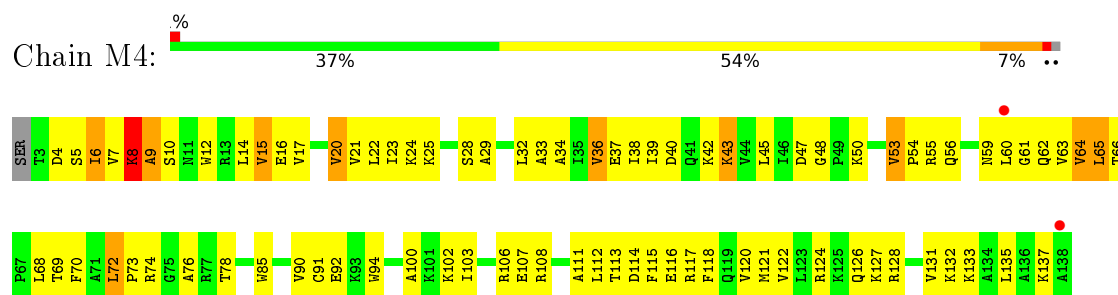


- Molecule 49: 60S ribosomal protein L13-A

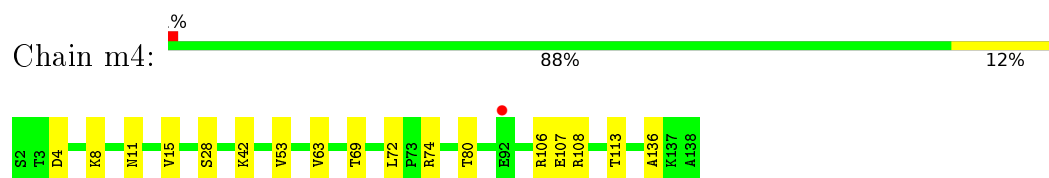




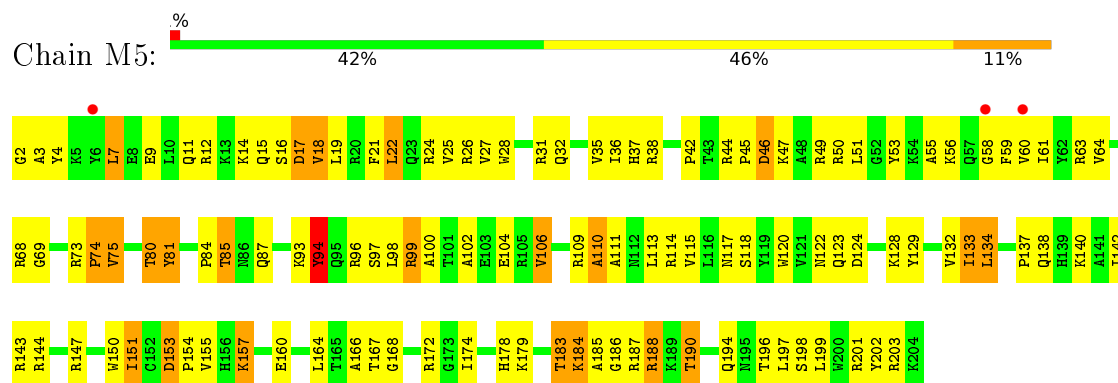
- Molecule 50: 60S ribosomal protein L14-A



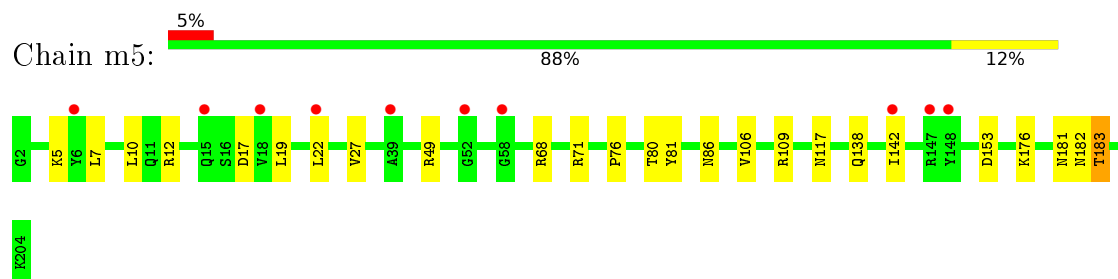
- Molecule 50: 60S ribosomal protein L14-A



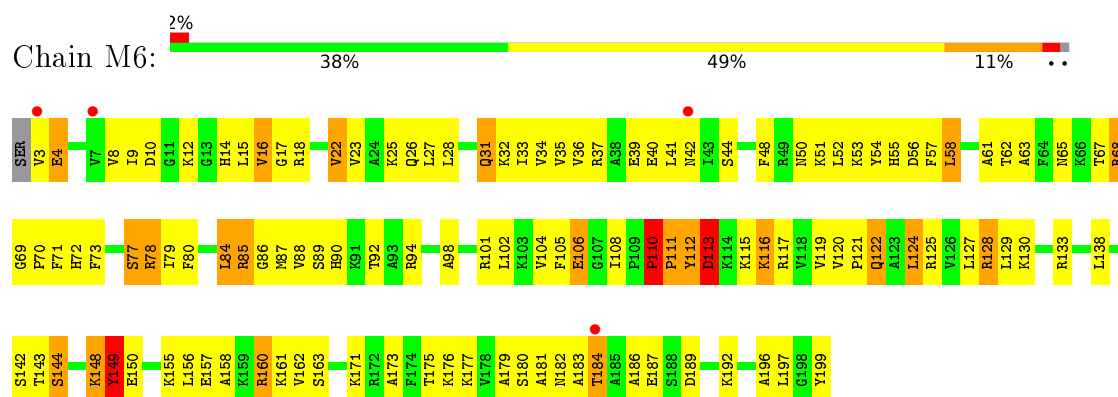
- Molecule 51: 60S ribosomal protein L15-A



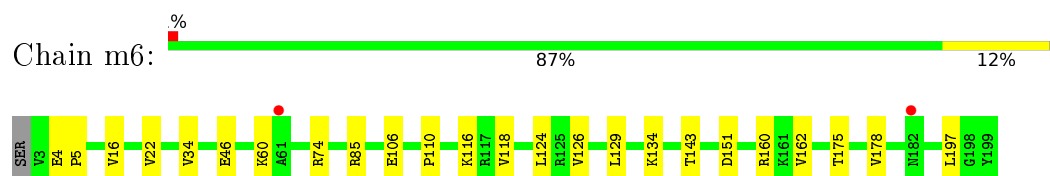
- Molecule 51: 60S ribosomal protein L15-A



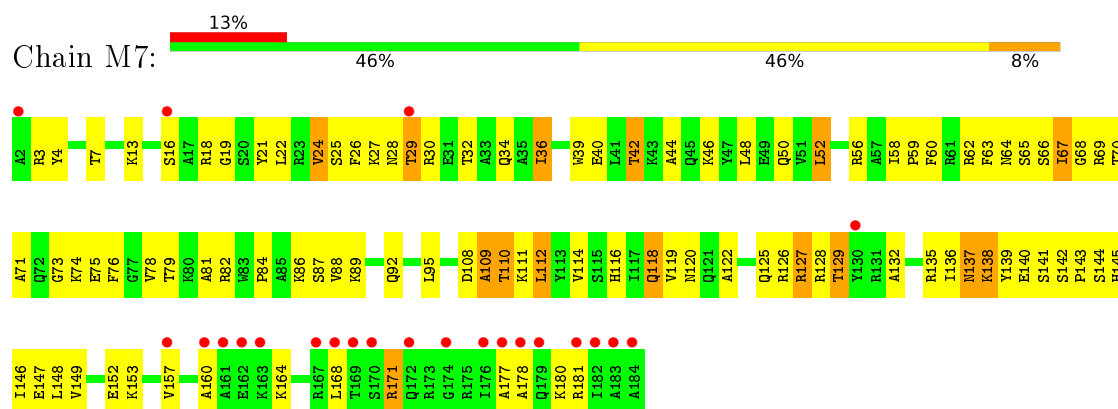
- Molecule 52: 60S ribosomal protein L16-A



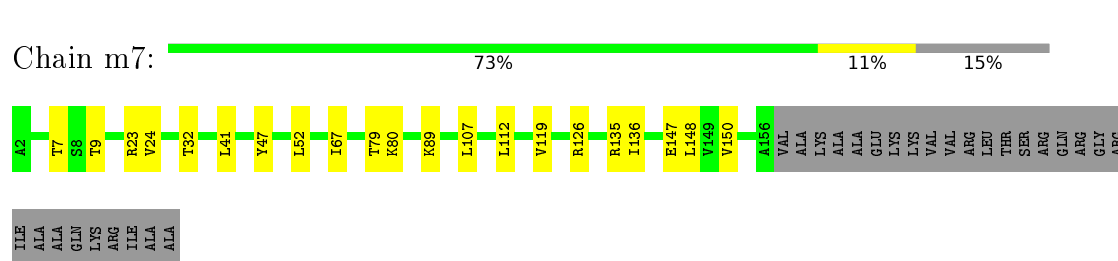
• Molecule 52: 60S ribosomal protein L16-A



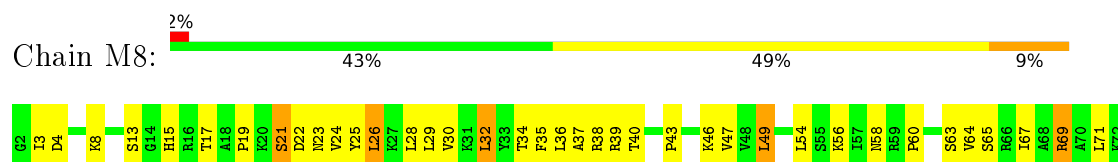
• Molecule 53: 60S ribosomal protein L17-A

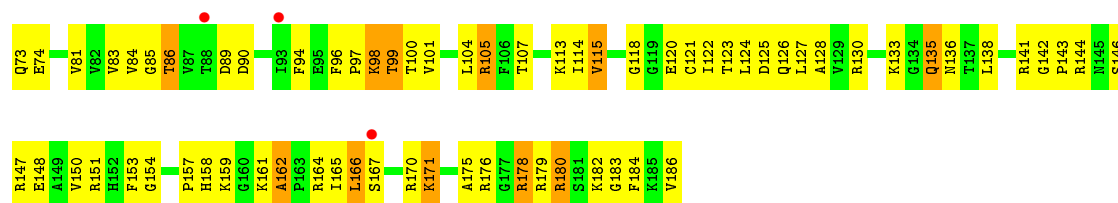


• Molecule 53: 60S ribosomal protein L17-A

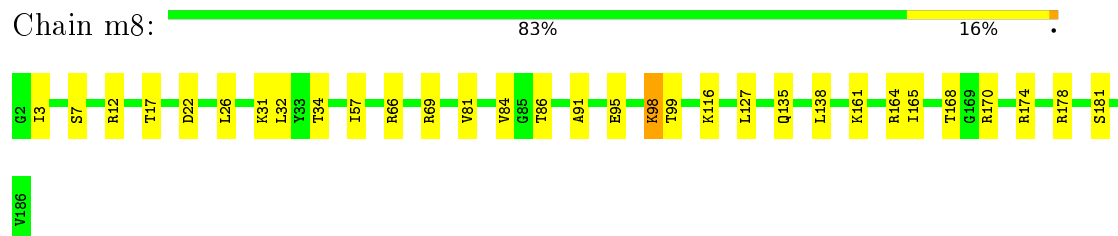


• Molecule 54: 60S ribosomal protein L18-A

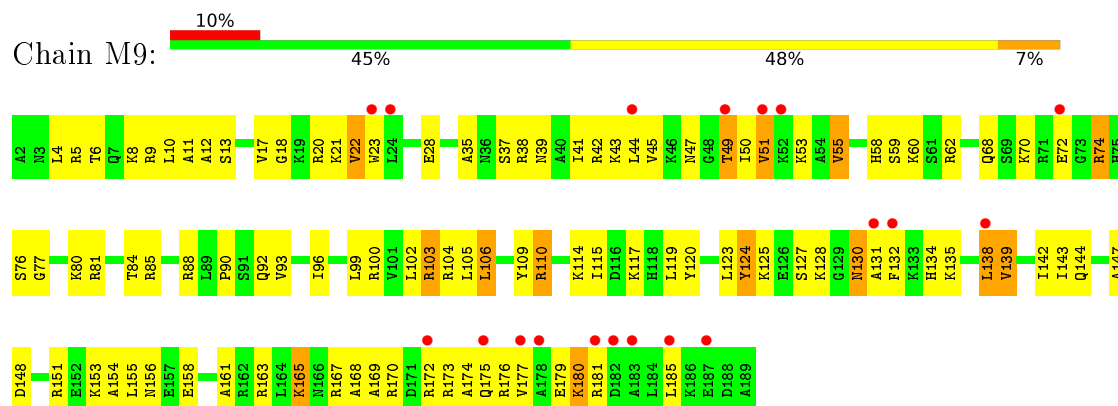




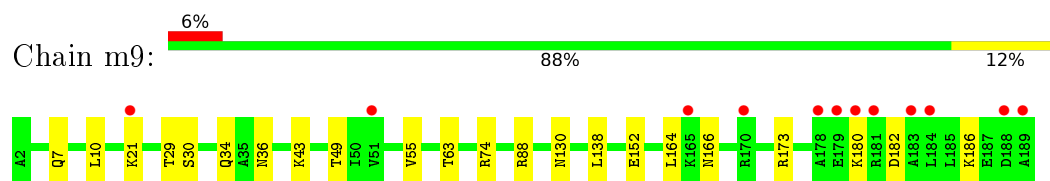
• Molecule 54: 60S ribosomal protein L18-A



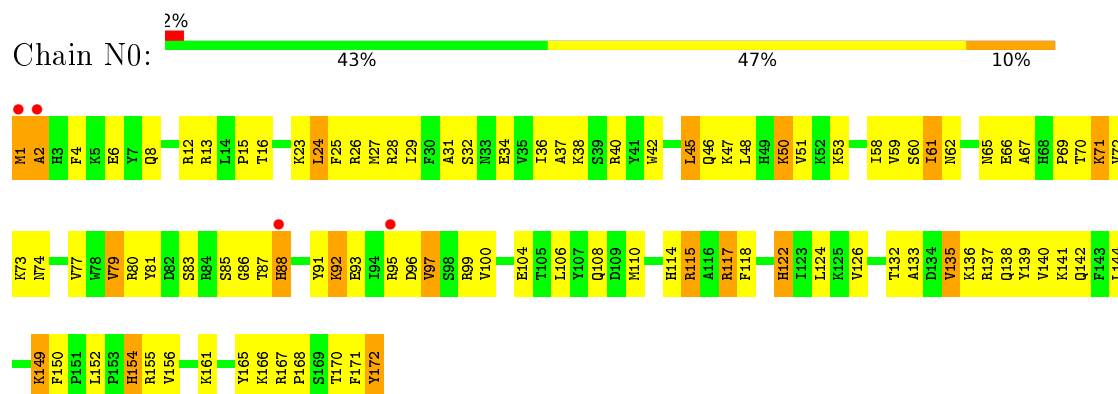
• Molecule 55: 60S ribosomal protein L19-A



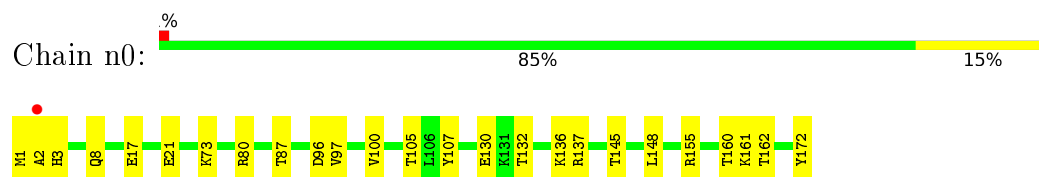
• Molecule 55: 60S ribosomal protein L19-A



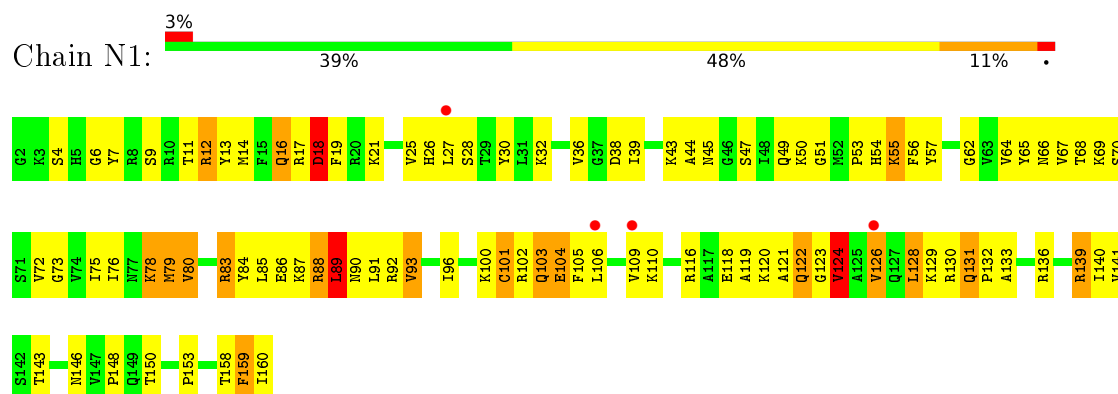
• Molecule 56: 60S ribosomal protein L20-A



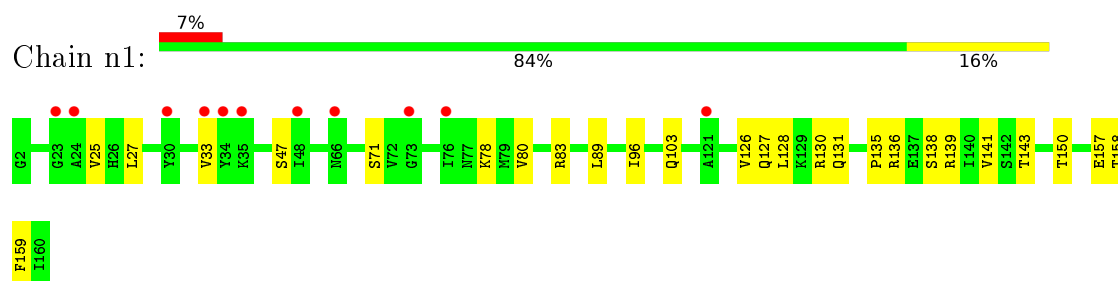
- Molecule 56: 60S ribosomal protein L20-A



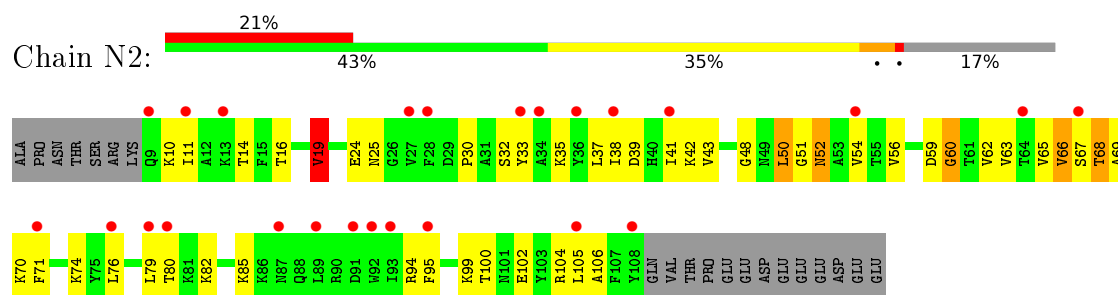
- Molecule 57: 60S ribosomal protein L21-A



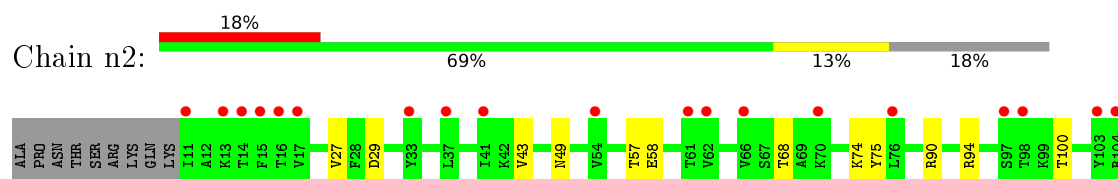
- Molecule 57: 60S ribosomal protein L21-A

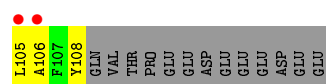


- Molecule 58: 60S ribosomal protein L22-A

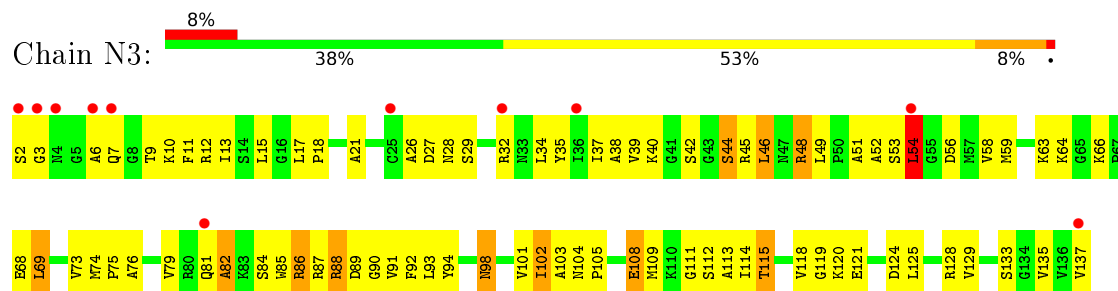


- Molecule 58: 60S ribosomal protein L22-A

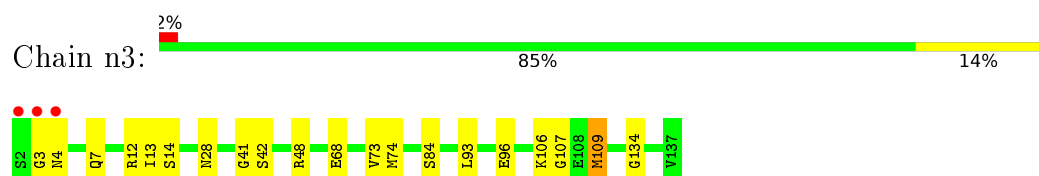




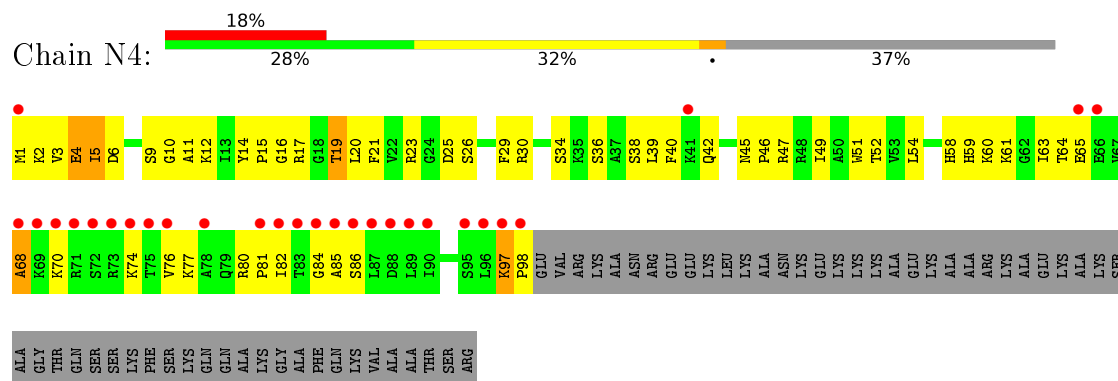
- Molecule 59: 60S ribosomal protein L23-A



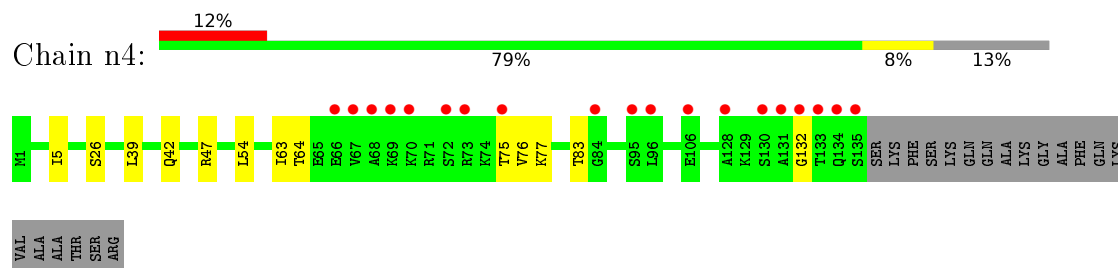
- Molecule 59: 60S ribosomal protein L23-A



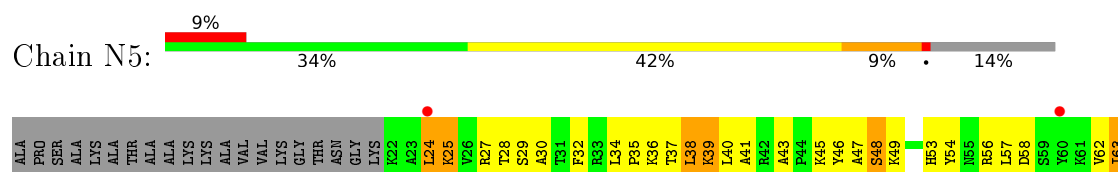
- Molecule 60: 60S ribosomal protein L24-A

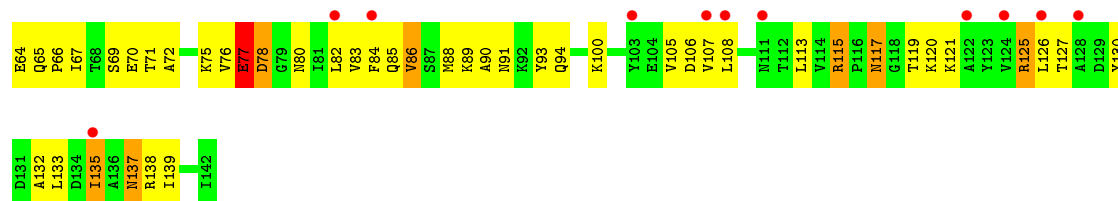


- Molecule 60: 60S ribosomal protein L24-A

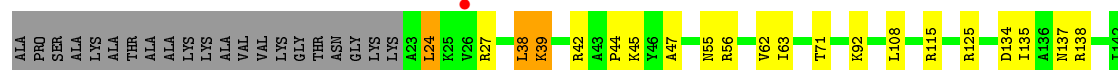


- Molecule 61: 60S ribosomal protein L25

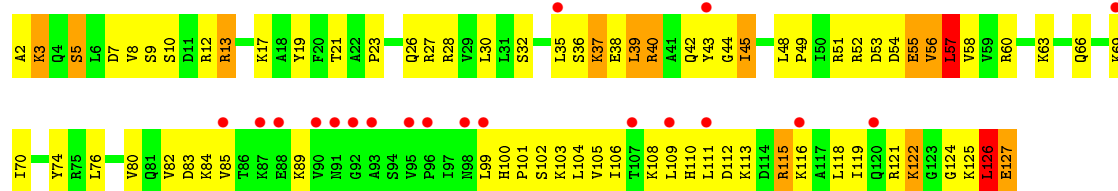
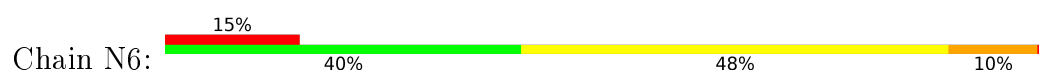




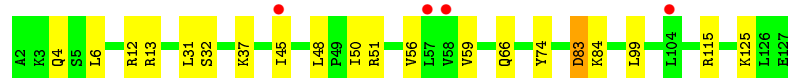
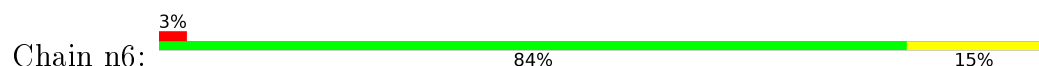
- Molecule 61: 60S ribosomal protein L25



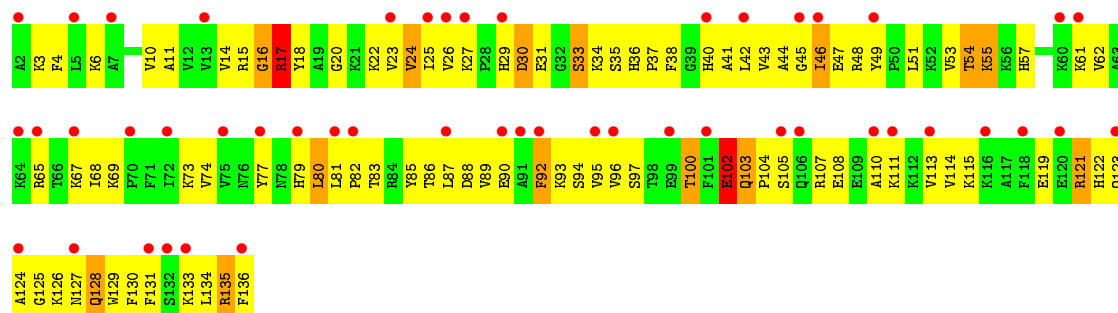
- Molecule 62: 60S ribosomal protein L26-A



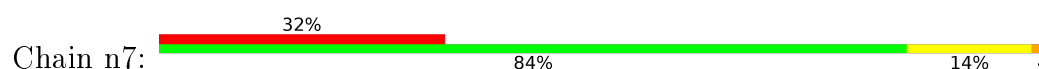
- Molecule 62: 60S ribosomal protein L26-A

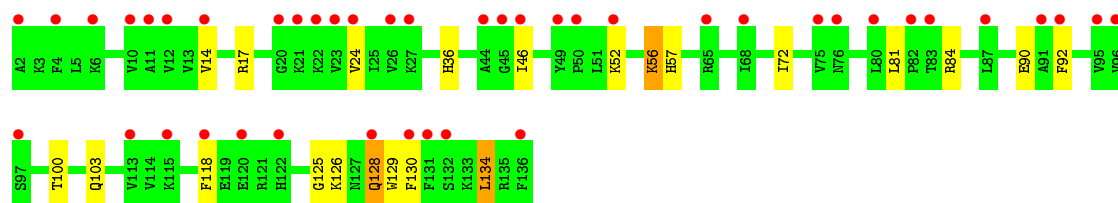


- Molecule 63: 60S ribosomal protein L27-A



- Molecule 63: 60S ribosomal protein L27-A

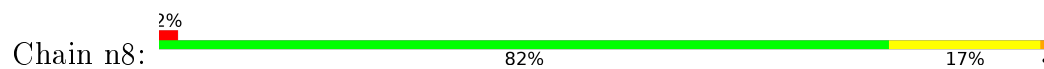




• Molecule 64: 60S ribosomal protein L28



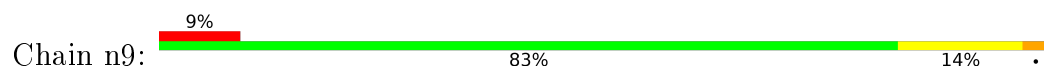
• Molecule 64: 60S ribosomal protein L28



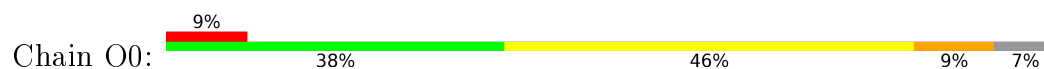
• Molecule 65: 60S ribosomal protein L29

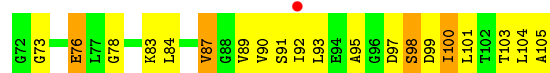


• Molecule 65: 60S ribosomal protein L29

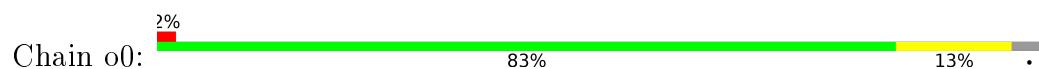


• Molecule 66: 60S ribosomal protein L30

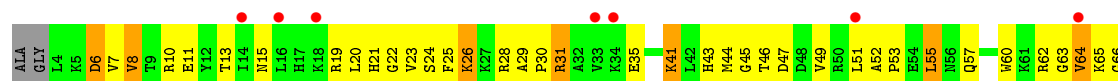
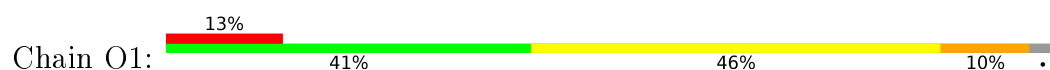




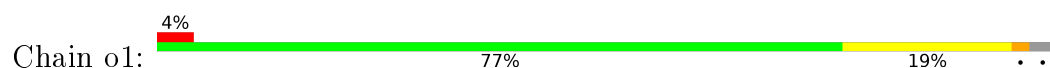
- Molecule 66: 60S ribosomal protein L30



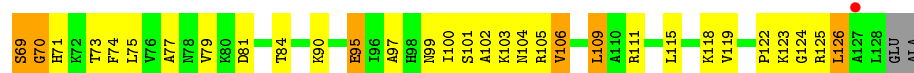
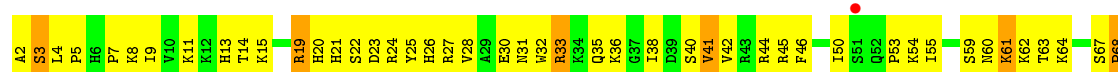
- Molecule 67: 60S ribosomal protein L31-A



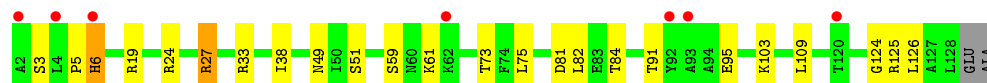
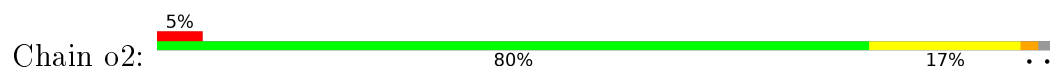
- Molecule 67: 60S ribosomal protein L31-A



- Molecule 68: 60S ribosomal protein L32



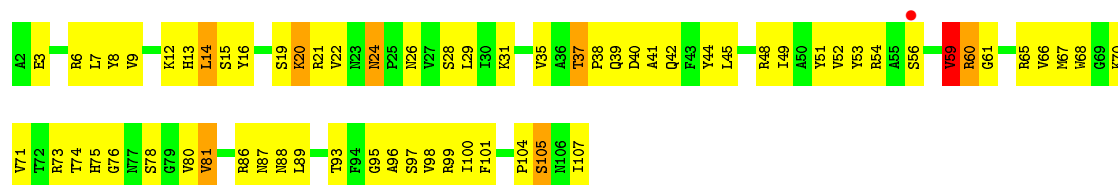
- Molecule 68: 60S ribosomal protein L32



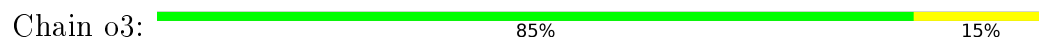
- Molecule 69: 60S ribosomal protein L33-A



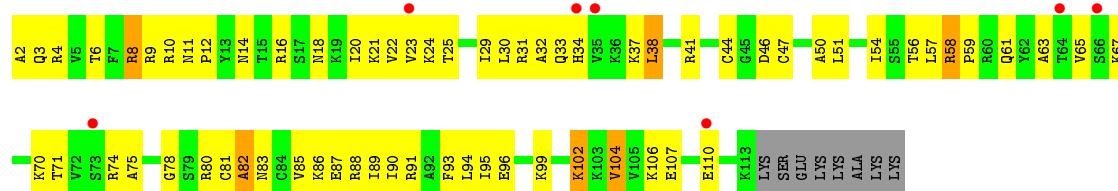
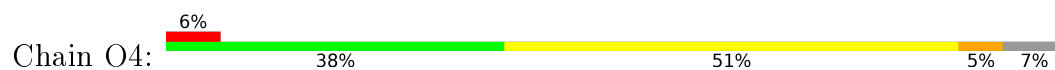




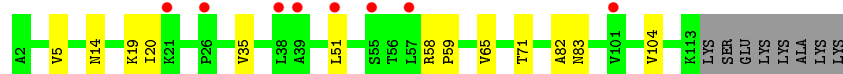
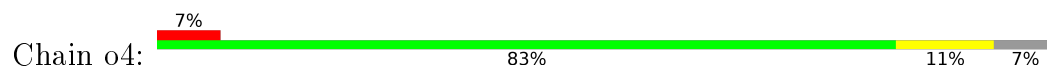
- Molecule 69: 60S ribosomal protein L33-A



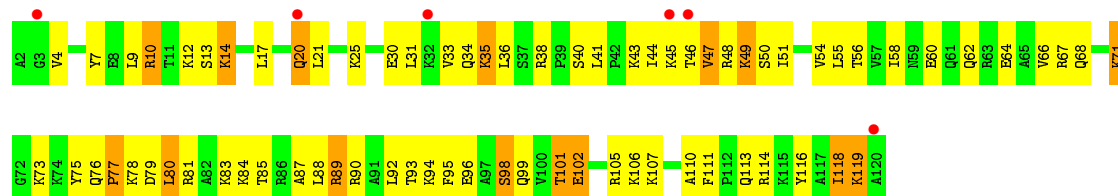
- Molecule 70: 60S ribosomal protein L34-A



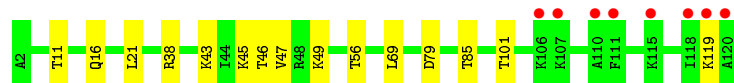
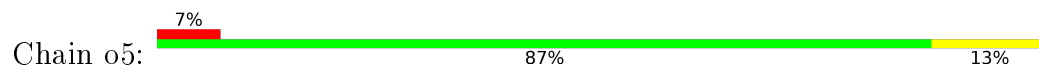
- Molecule 70: 60S ribosomal protein L34-A



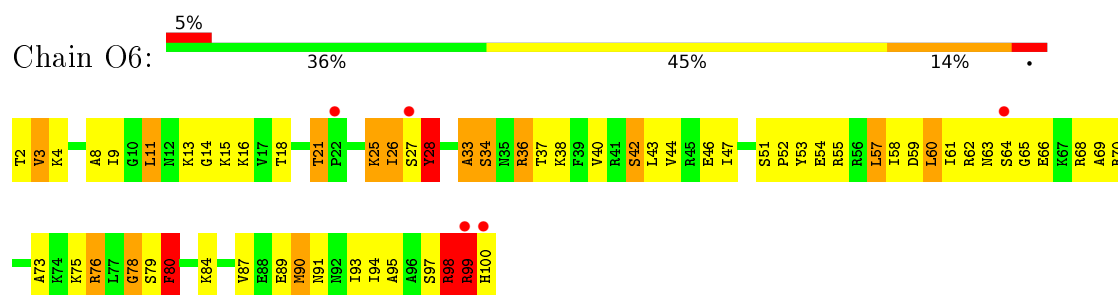
- Molecule 71: 60S ribosomal protein L35-A



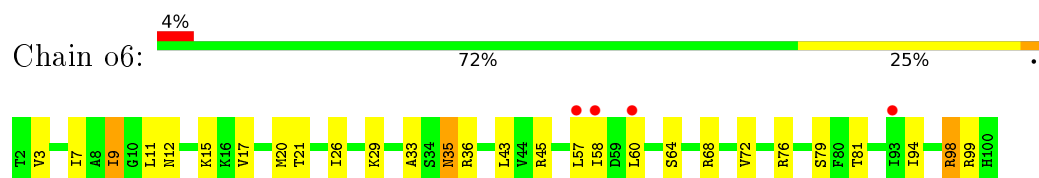
- Molecule 71: 60S ribosomal protein L35-A



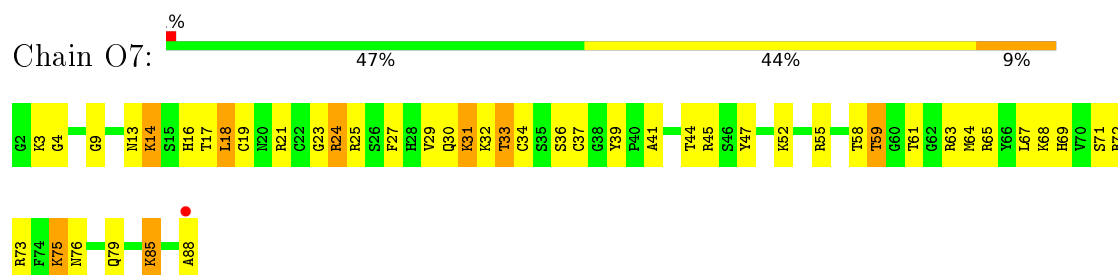
- Molecule 72: 60S ribosomal protein L36-A



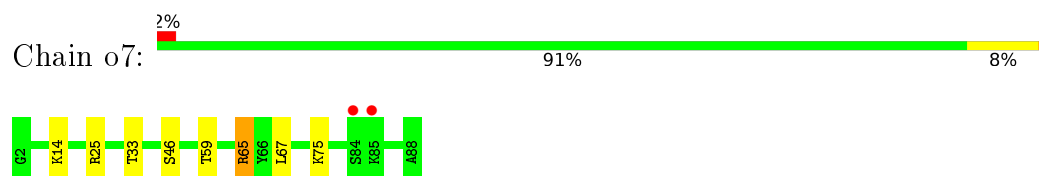
- Molecule 72: 60S ribosomal protein L36-A



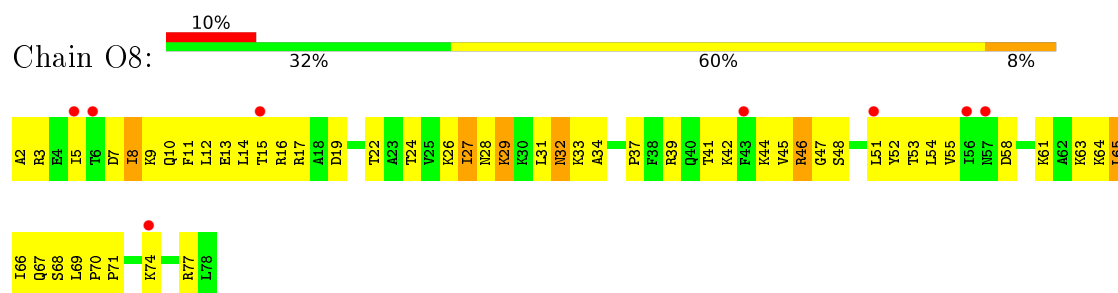
- Molecule 73: 60S ribosomal protein L37-A



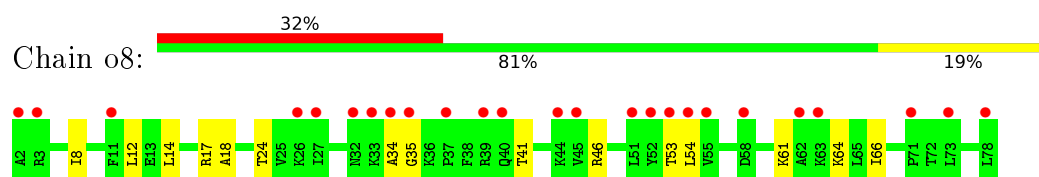
- Molecule 73: 60S ribosomal protein L37-A



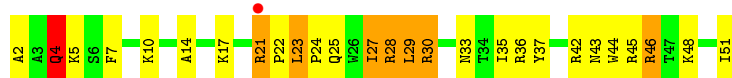
- Molecule 74: 60S ribosomal protein L38



- Molecule 74: 60S ribosomal protein L38



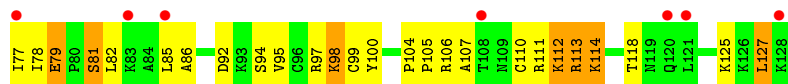
- Molecule 75: 60S ribosomal protein L39



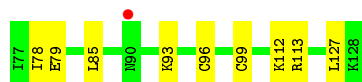
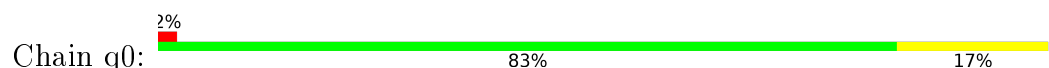
- Molecule 75: 60S ribosomal protein L39



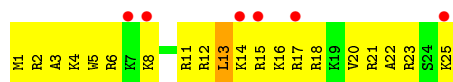
- Molecule 76: Ubiquitin-60S ribosomal protein L40



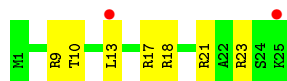
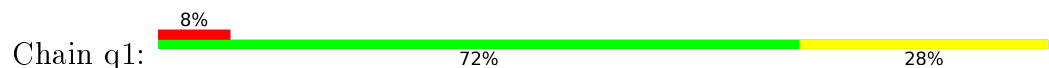
- Molecule 76: Ubiquitin-60S ribosomal protein L40



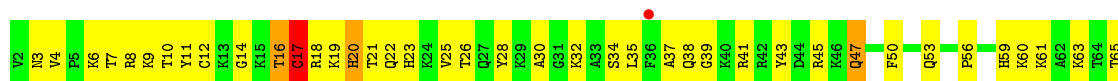
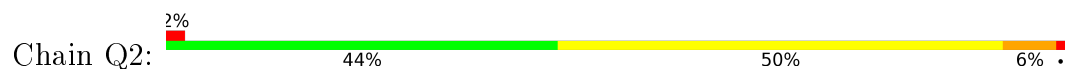
- Molecule 77: 60S ribosomal protein L41-A

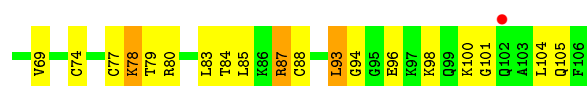


- Molecule 77: 60S ribosomal protein L41-A

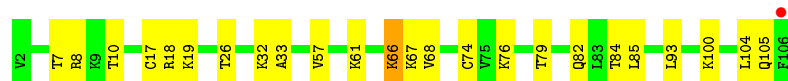
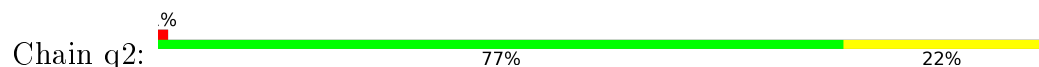


- Molecule 78: 60S ribosomal protein L42-A

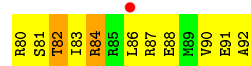
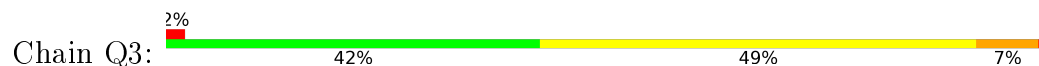




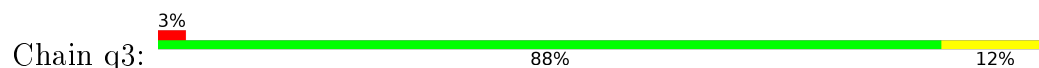
- Molecule 78: 60S ribosomal protein L42-A



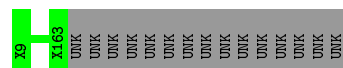
- Molecule 79: 60S ribosomal protein L43-A



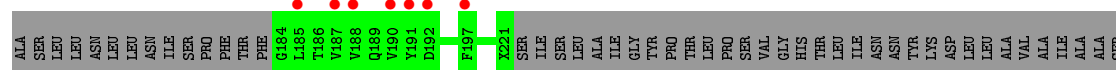
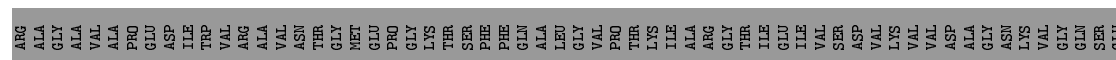
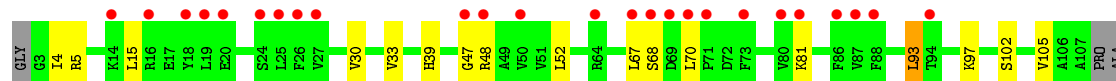
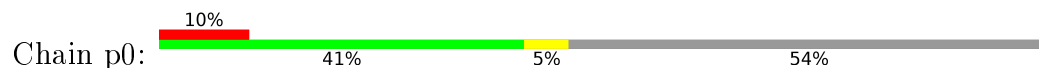
- Molecule 79: 60S ribosomal protein L43-A



- Molecule 80: 60S ribosomal protein L12-A (uL11)



- Molecule 81: 60S acidic ribosomal protein P0



TYR HIS  
TYR PRO  
GLU GLU  
ILE ASP  
LEU VAL  
ASP ARG  
ILE ILE  
GLU ASN  
PRO PRO  
GLU LYS  
TYR ALA  
ALA ALA  
ALA ALA  
ALA PRO  
ALA ALA  
ALA THR  
SER SER  
ALA ALA  
SER GLY  
ASP ASP  
ALA ALA  
ALA PRO  
ALA GLU  
GLU GLU  
ALA ALA  
ALA GLU  
GLU GLU  
GLU GLU  
SER SER  
ASP ASP  
ASP MET  
GLY PHE  
PHE LEU  
PHE ASP

- Molecule 82: 60S ribosomal protein P1 alpha/P2 beta

Chain p1:  100%

There are no outlier residues recorded for this chain.

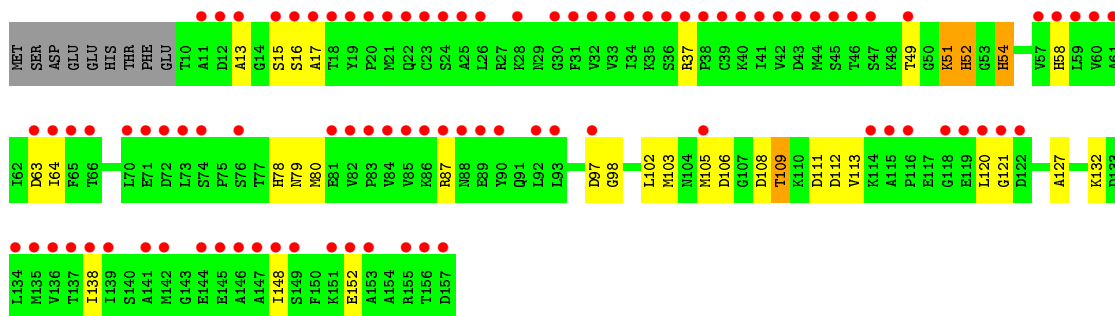
- Molecule 82: 60S ribosomal protein P1 alpha/P2 beta

Chain p2:  98%



- Molecule 83: Eukaryotic translation initiation factor 5A-1

Chain f:  59%  
73% 19% 6%



- Molecule 84: CH-CH-8AN-PRO-PRO

Chain B:  40% 60%



- Molecule 84: CH-CH-8AN-PRO-PRO

Chain C:  60% 40%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	435.58 Å   287.33 Å   303.82 Å 90.00°   98.95°   90.00°	Depositor
Resolution (Å)	99.53 – 3.30 99.53 – 3.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (99.53-3.30) 99.9 (99.53-3.30)	Depositor EDS
$R_{merge}$	0.59	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.34 (at 3.33 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.251 ,   0.310 0.258 ,   (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	85.9	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 74.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	404292	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CH, SPS, MG, OHX, 8AN

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	2	0.36	0/42467	0.89	42/66169 (0.1%)
1	6	0.43	0/42790	0.93	50/66673 (0.1%)
2	S0	0.28	0/1617	0.50	0/2215
2	s0	0.29	0/1623	0.51	0/2222
3	S1	0.26	0/1735	0.50	0/2335
3	s1	0.32	0/1748	0.53	0/2352
4	S2	0.30	0/1665	0.52	0/2263
4	s2	0.33	0/1665	0.55	1/2263 (0.0%)
5	S3	0.29	0/1759	0.47	0/2368
5	s3	0.27	0/1759	0.47	0/2368
6	S4	0.29	0/2109	0.53	0/2839
6	s4	0.34	0/2109	0.55	0/2839
7	S5	0.27	0/1629	0.50	0/2202
7	s5	0.28	0/1629	0.48	0/2202
8	S6	0.29	0/1823	0.47	0/2439
8	s6	0.33	0/1779	0.53	0/2379
9	S7	0.28	0/1506	0.53	1/2028 (0.0%)
9	s7	0.29	0/1516	0.52	0/2043
10	S8	0.30	0/1514	0.51	0/2021
10	s8	0.34	0/1514	0.52	0/2021
11	S9	0.29	0/1519	0.48	0/2035
11	s9	0.31	0/1519	0.50	0/2035
12	C0	0.28	0/725	0.44	0/978
12	c0	0.27	0/713	0.51	1/961 (0.1%)
13	C1	0.31	0/1185	0.50	0/1598
13	c1	0.36	0/1185	0.55	0/1598
14	C2	0.26	0/819	0.49	0/1109
14	c2	0.22	0/819	0.47	0/1109
15	C3	0.31	0/1215	0.52	1/1638 (0.1%)
15	c3	0.33	0/1215	0.51	0/1638
16	C4	0.26	0/901	0.51	0/1217
16	c4	0.30	0/960	0.54	0/1290

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
17	C5	0.31	0/988	0.54	0/1327
17	c5	0.29	0/1030	0.52	0/1384
18	C6	0.29	0/1125	0.54	2/1510 (0.1%)
18	c6	0.29	0/1131	0.51	0/1518
19	C7	0.30	0/935	0.52	0/1254
19	c7	0.27	0/914	0.50	0/1224
20	C8	0.30	0/1211	0.51	0/1628
20	c8	0.28	0/1211	0.49	0/1628
21	C9	0.29	0/1130	0.48	0/1517
21	c9	0.29	0/1130	0.47	0/1517
22	D0	0.30	0/865	0.52	0/1169
22	d0	0.29	0/892	0.49	0/1205
23	D1	0.29	0/693	0.49	0/935
23	d1	0.31	0/693	0.50	0/935
24	D2	0.30	0/1038	0.55	1/1395 (0.1%)
24	d2	0.35	0/1038	0.56	1/1395 (0.1%)
25	D3	0.35	0/1139	0.55	0/1518
25	d3	0.38	0/1139	0.57	0/1518
26	D4	0.29	0/1087	0.47	0/1449
26	d4	0.31	0/1087	0.51	0/1449
27	D5	0.26	0/571	0.55	0/768
27	d5	0.26	0/566	0.47	0/761
28	D6	0.29	0/782	0.53	0/1047
28	d6	0.34	0/782	0.53	0/1047
29	D7	0.29	0/620	0.51	0/838
29	d7	0.28	0/620	0.49	0/838
30	D8	0.27	0/499	0.47	0/670
30	d8	0.28	0/499	0.49	0/670
31	D9	0.32	0/452	0.55	1/600 (0.2%)
31	d9	0.33	0/452	0.50	0/600
32	E0	0.28	0/483	0.48	0/643
32	e0	0.32	0/499	0.54	0/665
33	E1	0.29	0/577	0.56	0/770
33	e1	0.29	0/619	0.58	0/822
34	SR	0.26	0/2490	0.49	0/3389
34	sR	0.25	0/2495	0.44	0/3395
35	SM	0.31	0/925	0.54	1/1240 (0.1%)
35	sM	0.31	0/480	0.53	0/642
36	1	0.56	0/75394	1.03	115/117545 (0.1%)
36	5	0.59	2/75865 (0.0%)	1.05	143/118275 (0.1%)
37	3	0.46	0/2883	0.91	0/4491
37	7	0.55	0/2883	1.00	2/4491 (0.0%)
38	4	0.53	0/3746	1.00	5/5832 (0.1%)



Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
38	8	0.52	0/3746	0.97	2/5832 (0.0%)
39	L2	0.37	0/1948	0.57	0/2617
39	l2	0.40	0/1946	0.62	0/2614
40	L3	0.39	0/3146	0.56	0/4228
40	l3	0.45	0/3146	0.58	0/4228
41	L4	0.41	0/2800	0.61	0/3790
41	l4	0.41	1/2800 (0.0%)	0.63	2/3790 (0.1%)
42	L5	0.35	0/2425	0.53	0/3271
42	l5	0.42	0/2408	0.55	0/3248
43	L6	0.39	0/1260	0.55	0/1694
43	l6	0.41	0/1269	0.57	0/1705
44	L7	0.42	0/1821	0.59	0/2451
44	l7	0.43	0/1828	0.62	2/2461 (0.1%)
45	L8	0.31	0/1836	0.53	0/2481
45	l8	0.32	0/1796	0.51	0/2431
46	L9	0.36	0/1539	0.56	0/2073
46	l9	0.40	0/1539	0.57	0/2073
47	M0	0.40	1/1741 (0.1%)	0.55	0/2335
47	m0	0.43	0/1758	0.61	2/2358 (0.1%)
48	M1	0.31	0/1374	0.50	0/1842
48	m1	0.38	0/1374	0.59	1/1842 (0.1%)
49	M3	0.40	0/1568	0.59	0/2106
49	m3	0.37	0/1573	0.57	0/2113
50	M4	0.40	0/1068	0.54	0/1438
50	m4	0.41	0/1074	0.57	0/1446
51	M5	0.38	0/1757	0.56	0/2354
51	m5	0.36	0/1757	0.55	0/2354
52	M6	0.44	0/1585	0.59	0/2128
52	m6	0.50	0/1585	0.61	0/2128
53	M7	0.41	0/1443	0.58	0/1944
53	m7	0.44	0/1250	0.59	0/1683
54	M8	0.39	0/1465	0.60	0/1965
54	m8	0.39	0/1465	0.60	0/1965
55	M9	0.31	0/1538	0.49	0/2050
55	m9	0.34	0/1538	0.51	0/2050
56	N0	0.39	0/1481	0.56	0/1990
56	n0	0.44	0/1481	0.59	0/1990
57	N1	0.40	0/1300	0.56	1/1743 (0.1%)
57	n1	0.43	0/1300	0.55	0/1743
58	N2	0.30	0/812	0.48	0/1099
58	n2	0.32	0/794	0.52	0/1076
59	N3	0.39	0/1018	0.57	0/1369
59	n3	0.45	0/1018	0.61	0/1369

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
60	N4	0.31	0/712	0.50	0/958
60	n4	0.35	0/1052	0.51	0/1398
61	N5	0.33	0/979	0.55	0/1321
61	n5	0.36	0/974	0.57	0/1314
62	N6	0.35	0/1004	0.62	2/1341 (0.1%)
62	n6	0.35	0/1004	0.57	0/1341
63	N7	0.30	0/1118	0.50	0/1497
63	n7	0.31	0/1118	0.51	0/1497
64	N8	0.41	0/1204	0.62	0/1612
64	n8	0.40	0/1204	0.61	0/1612
65	N9	0.38	0/473	0.60	0/629
65	n9	0.43	0/473	0.67	0/629
66	O0	0.30	0/751	0.47	0/1008
66	o0	0.31	0/775	0.52	0/1040
67	O1	0.34	0/890	0.52	0/1196
67	o1	0.41	0/897	0.60	0/1205
68	O2	0.43	0/1041	0.63	0/1394
68	o2	0.44	0/1041	0.62	0/1394
69	O3	0.48	0/868	0.56	0/1168
69	o3	0.48	0/868	0.58	0/1168
70	O4	0.34	0/890	0.54	0/1189
70	o4	0.34	0/890	0.58	0/1189
71	O5	0.37	0/978	0.57	0/1301
71	o5	0.34	0/974	0.53	0/1297
72	O6	0.37	0/778	0.60	0/1034
72	o6	0.34	0/777	0.54	0/1033
73	O7	0.42	0/696	0.61	0/923
73	o7	0.40	0/696	0.59	0/923
74	O8	0.30	0/618	0.47	0/826
74	o8	0.31	0/614	0.54	0/822
75	O9	0.39	0/443	0.58	0/588
75	o9	0.39	0/443	0.53	0/588
76	Q0	0.43	0/423	0.60	0/562
76	q0	0.45	0/423	0.55	0/562
77	Q1	0.36	0/234	0.59	0/300
77	q1	0.43	0/234	0.62	0/300
78	Q2	0.53	1/860 (0.1%)	0.63	0/1136
78	q2	0.52	1/860 (0.1%)	0.63	1/1136 (0.1%)
79	Q3	0.37	0/701	0.59	0/934
79	q3	0.43	0/701	0.56	0/934
81	p0	0.27	0/977	0.48	0/1313
83	f	0.34	0/1131	0.58	0/1522
84	B	0.31	0/15	0.54	0/20

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
84	C	0.41	0/15	0.80	0/20
All	All	0.46	6/431576 (0.0%)	0.85	380/633602 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
18	c6	0	1
19	C7	0	1
27	D5	0	1
39	l2	0	1
44	L7	0	1
52	M6	0	1
64	n8	0	1
65	N9	0	1
65	n9	0	1
83	f	0	1
All	All	0	10

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
78	q2	17	CYS	CB-SG	8.77	1.97	1.82
78	Q2	17	CYS	CB-SG	7.99	1.95	1.82
36	5	1152	G	N9-C4	-5.86	1.33	1.38
36	5	942	U	C4-O4	5.25	1.27	1.23
47	M0	8	CYS	CB-SG	-5.03	1.73	1.81
41	14	94	CYS	CB-SG	-5.00	1.73	1.81

All (380) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	1152	G	N3-C4-N9	-11.60	119.04	126.00
36	5	1152	G	N3-C4-C5	11.31	134.25	128.60
36	5	2726	C	C6-N1-C2	-9.68	116.43	120.30
36	5	1152	G	C2-N3-C4	-9.61	107.10	111.90
36	1	406	G	O4'-C1'-N9	9.54	115.83	108.20
36	1	1308	A	C8-N9-C4	-9.32	102.07	105.80
36	1	2617	U	C5-C4-O4	8.97	131.28	125.90
36	5	2278	C	N1-C2-O2	8.83	124.20	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
38	4	125	U	C2-N1-C1'	8.67	128.11	117.70
36	1	1308	A	N7-C8-N9	8.59	118.09	113.80
36	5	2169	G	C5-C6-O6	8.39	133.64	128.60
36	1	3217	C	C2-N1-C1'	8.35	127.98	118.80
1	6	453	U	C2-N1-C1'	8.21	127.56	117.70
36	5	1028	U	C2-N1-C1'	8.14	127.46	117.70
36	1	3217	C	N1-C2-O2	8.13	123.78	118.90
36	1	439	C	N1-C2-O2	8.12	123.77	118.90
36	1	2617	U	N1-C2-N3	8.03	119.72	114.90
36	5	2726	C	N3-C2-O2	-7.94	116.34	121.90
36	1	2355	G	N1-C6-O6	7.70	124.52	119.90
36	5	942	U	N3-C4-C5	-7.70	109.98	114.60
36	5	2524	A	O4'-C1'-N9	7.62	114.30	108.20
1	6	163	G	N3-C4-N9	-7.59	121.45	126.00
36	1	1115	G	C6-C5-N7	-7.56	125.86	130.40
36	5	1200	A	N1-C6-N6	7.56	123.13	118.60
36	5	1192	C	C5-C6-N1	7.53	124.76	121.00
36	1	2403	G	N1-C6-O6	7.51	124.41	119.90
36	1	2617	U	C4-C5-C6	7.38	124.13	119.70
36	1	3217	C	N3-C2-O2	-7.38	116.74	121.90
36	5	1152	G	C8-N9-C1'	7.37	136.58	127.00
1	6	1473	U	C2-N1-C1'	7.35	126.52	117.70
36	5	2169	G	C8-N9-C4	-7.33	103.47	106.40
36	5	2169	G	C4-C5-N7	-7.32	107.87	110.80
36	1	439	C	C2-N1-C1'	7.32	126.85	118.80
1	6	453	U	N1-C2-O2	7.29	127.90	122.80
36	5	3351	U	C2-N1-C1'	7.28	126.44	117.70
1	6	1473	U	N1-C2-O2	7.28	127.90	122.80
36	1	1367	G	N1-C6-O6	7.27	124.26	119.90
1	2	934	C	C2-N1-C1'	7.27	126.79	118.80
36	1	2714	G	N3-C4-C5	7.24	132.22	128.60
62	N6	57	LEU	CA-CB-CG	7.23	131.93	115.30
36	5	2726	C	C2-N1-C1'	7.23	126.75	118.80
36	5	1028	U	N1-C2-O2	7.21	127.84	122.80
36	5	1495	U	C2-N1-C1'	7.18	126.32	117.70
36	1	3278	C	N1-C2-O2	7.16	123.20	118.90
36	1	1115	G	N3-C4-N9	7.16	130.29	126.00
36	1	2869	U	C5-C4-O4	-7.14	121.62	125.90
36	5	406	G	N1-C6-O6	-7.14	115.62	119.90
1	2	831	U	C2-N1-C1'	7.14	126.27	117.70
36	5	1152	G	C4-N9-C1'	-7.08	117.30	126.50
1	6	1000	C	C2-N1-C1'	7.04	126.55	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
44	17	229	PHE	CB-CG-CD1	7.03	125.72	120.80
36	5	2272	G	O4'-C1'-N9	7.03	113.82	108.20
36	1	1556	C	N1-C2-O2	6.97	123.08	118.90
36	5	568	G	N1-C6-O6	6.96	124.08	119.90
36	5	2827	U	N1-C2-O2	6.95	127.67	122.80
1	2	1698	G	P-O3'-C3'	6.91	127.99	119.70
1	6	1274	C	N1-C2-O2	6.91	123.05	118.90
36	1	3278	C	N3-C2-O2	-6.91	117.06	121.90
1	6	680	U	C2-N1-C1'	6.91	125.99	117.70
36	5	2283	G	N1-C6-O6	6.88	124.03	119.90
36	1	3057	U	N3-C2-O2	-6.84	117.41	122.20
36	1	656	A	N1-C6-N6	6.83	122.70	118.60
1	2	728	U	N1-C2-O2	6.83	127.58	122.80
36	5	419	G	N1-C6-O6	-6.80	115.82	119.90
1	6	1473	U	N3-C2-O2	-6.76	117.47	122.20
1	2	1389	C	N1-C2-O2	6.76	122.96	118.90
36	5	1308	A	O5'-P-OP1	-6.71	99.66	105.70
38	4	125	U	N1-C2-O2	6.68	127.48	122.80
36	1	2819	A	O5'-P-OP2	-6.68	99.69	105.70
36	5	343	U	O5'-P-OP2	-6.65	99.71	105.70
1	6	680	U	N1-C2-O2	6.63	127.44	122.80
1	6	813	U	C2-N1-C1'	6.59	125.61	117.70
1	2	934	C	N1-C2-O2	6.57	122.84	118.90
36	5	2938	G	N1-C6-O6	-6.54	115.97	119.90
1	2	959	U	C2-N1-C1'	6.52	125.52	117.70
1	6	813	U	N1-C2-O2	6.47	127.33	122.80
36	5	2211	U	N3-C2-O2	-6.46	117.68	122.20
36	5	880	G	C4-N9-C1'	-6.45	118.12	126.50
36	5	568	G	C5-C6-N1	-6.45	108.28	111.50
1	2	728	U	C2-N1-C1'	6.44	125.42	117.70
36	1	1367	G	C5-C6-N1	-6.44	108.28	111.50
36	1	835	G	O4'-C1'-N9	6.42	113.34	108.20
36	5	3154	C	N1-C2-O2	6.42	122.75	118.90
1	2	728	U	N3-C2-O2	-6.42	117.71	122.20
1	2	831	U	C5-C6-N1	6.40	125.90	122.70
1	6	1274	C	C2-N1-C1'	6.40	125.84	118.80
36	1	1556	C	N3-C2-O2	-6.39	117.43	121.90
36	5	2372	A	P-O3'-C3'	6.37	127.34	119.70
36	1	650	C	N3-C2-O2	6.37	126.36	121.90
36	5	1367	G	N1-C6-O6	6.36	123.72	119.90
36	5	1307	G	P-O3'-C3'	6.35	127.32	119.70
36	5	1200	A	N9-C4-C5	-6.34	103.26	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	2572	C	C2-N1-C1'	6.33	125.77	118.80
36	5	2957	G	O5'-P-OP1	-6.32	100.01	105.70
36	1	1115	G	C4-N9-C1'	6.32	134.71	126.50
36	5	2169	G	N9-C4-C5	6.31	107.92	105.40
1	2	1389	C	C2-N1-C1'	6.30	125.73	118.80
1	6	1097	U	P-O3'-C3'	6.30	127.26	119.70
1	6	453	U	N3-C2-O2	-6.28	117.81	122.20
36	5	2847	A	C8-N9-C4	6.27	108.31	105.80
18	C6	40	GLU	C-N-CD	-6.26	106.83	120.60
38	4	125	U	C6-N1-C1'	-6.25	112.44	121.20
36	5	2860	U	O4'-C1'-N1	6.25	113.20	108.20
36	5	645	A	N1-C6-N6	-6.25	114.85	118.60
1	6	813	U	N3-C2-O2	-6.24	117.83	122.20
36	5	406	G	C8-N9-C1'	6.23	135.10	127.00
1	2	1052	U	C2-N1-C1'	6.22	125.16	117.70
36	5	1028	U	N3-C2-O2	-6.22	117.85	122.20
12	c0	83	PRO	N-CA-CB	6.21	110.75	103.30
36	1	2808	A	N1-C6-N6	6.17	122.31	118.60
36	1	1367	G	C6-C5-N7	-6.16	126.70	130.40
36	5	2572	C	N1-C2-O2	6.15	122.59	118.90
36	5	1352	A	OP1-P-O3'	6.14	118.71	105.20
36	1	2403	G	C6-C5-N7	-6.14	126.72	130.40
1	6	647	G	N3-C4-N9	-6.13	122.32	126.00
36	5	2234	G	N1-C6-O6	6.12	123.58	119.90
36	5	641	C	C2-N1-C1'	-6.12	112.07	118.80
36	1	583	G	O5'-P-OP1	-6.09	100.22	105.70
36	1	1115	G	C8-N9-C1'	-6.08	119.10	127.00
36	1	922	U	C2-N1-C1'	6.06	124.97	117.70
1	2	959	U	N3-C2-O2	-6.05	117.96	122.20
36	1	2374	C	C2-N1-C1'	6.01	125.41	118.80
1	6	275	C	C2-N1-C1'	6.00	125.40	118.80
1	2	959	U	N1-C2-O2	5.99	126.99	122.80
37	7	73	C	C6-N1-C2	-5.98	117.91	120.30
36	1	98	G	C6-C5-N7	-5.98	126.81	130.40
1	6	680	U	N3-C2-O2	-5.97	118.02	122.20
36	5	1302	A	O5'-P-OP1	-5.97	100.33	105.70
36	1	650	C	N1-C2-O2	-5.96	115.32	118.90
1	2	158	U	P-O3'-C3'	5.96	126.85	119.70
36	1	406	G	C4-N9-C1'	-5.96	118.75	126.50
36	5	2860	U	C5-C4-O4	5.96	129.47	125.90
36	5	1113	G	C8-N9-C4	-5.96	104.02	106.40
1	6	937	C	C6-N1-C2	-5.95	117.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	1115	G	N3-C4-N9	5.95	129.57	126.00
36	1	3095	U	C5-C6-N1	5.94	125.67	122.70
36	1	282	G	C8-N9-C4	-5.93	104.03	106.40
36	5	1208	U	N3-C2-O2	-5.92	118.05	122.20
36	1	2836	C	N3-C2-O2	-5.91	117.76	121.90
36	1	3217	C	C6-N1-C1'	-5.90	113.72	120.80
36	1	2627	C	N1-C2-O2	-5.90	115.36	118.90
1	2	1456	C	C2-N1-C1'	5.90	125.29	118.80
1	6	542	A	P-O3'-C3'	5.90	126.78	119.70
1	6	280	U	C2-N1-C1'	5.89	124.77	117.70
36	5	1177	G	N1-C6-O6	5.89	123.44	119.90
36	5	3245	A	N1-C6-N6	5.88	122.13	118.60
36	5	880	G	C8-N9-C1'	5.87	134.63	127.00
36	5	406	G	C4-N9-C1'	-5.87	118.87	126.50
36	1	2608	G	N1-C6-O6	5.86	123.42	119.90
36	5	1481	A	P-O3'-C3'	5.86	126.73	119.70
1	6	1185	U	N1-C2-O2	5.84	126.89	122.80
36	1	3275	U	C2-N1-C1'	5.83	124.70	117.70
36	5	1166	G	N7-C8-N9	-5.83	110.19	113.10
36	5	2620	G	N3-C4-N9	-5.83	122.50	126.00
37	7	92	A	C8-N9-C4	5.83	108.13	105.80
36	5	1433	A	C8-N9-C4	-5.83	103.47	105.80
36	5	3245	A	N7-C8-N9	5.82	116.71	113.80
36	5	3154	C	C2-N1-C1'	5.82	125.20	118.80
36	5	835	G	O4'-C1'-N9	5.81	112.85	108.20
41	14	339	LEU	CA-CB-CG	5.80	128.64	115.30
36	5	3245	A	C6-C5-N7	-5.79	128.25	132.30
1	2	1389	C	N3-C2-O2	-5.78	117.85	121.90
36	1	2187	G	C5-C6-O6	5.78	132.07	128.60
38	8	80	A	N7-C8-N9	5.78	116.69	113.80
1	6	1389	C	C2-N1-C1'	5.77	125.15	118.80
36	1	2983	C	O4'-C1'-N1	5.77	112.82	108.20
38	4	39	G	N1-C6-O6	-5.77	116.44	119.90
1	6	1306	C	C2-N1-C1'	5.76	125.14	118.80
36	1	1115	G	C4-C5-N7	5.76	113.10	110.80
36	1	2873	U	C5-C4-O4	5.76	129.36	125.90
36	1	439	C	N3-C2-O2	-5.75	117.87	121.90
38	8	80	A	C8-N9-C4	-5.75	103.50	105.80
1	2	1202	A	C8-N9-C4	-5.75	103.50	105.80
36	5	2142	A	C8-N9-C4	5.74	108.10	105.80
36	5	406	G	C6-C5-N7	5.74	133.84	130.40
36	5	1887	A	O5'-P-OP2	-5.73	100.54	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	2231	C	C2-N1-C1'	5.73	125.10	118.80
36	5	1367	G	C5-C6-N1	-5.73	108.64	111.50
36	5	2526	C	N1-C2-O2	5.72	122.33	118.90
36	1	2403	G	C4-C5-C6	5.70	122.22	118.80
36	1	2617	U	N3-C4-C5	-5.69	111.19	114.60
36	1	637	C	C6-N1-C2	5.68	122.57	120.30
24	d2	93	LEU	CA-CB-CG	5.67	128.35	115.30
1	6	321	C	C2-N1-C1'	5.66	125.03	118.80
36	1	644	G	C5-C6-O6	5.66	132.00	128.60
36	5	1166	G	C8-N9-C4	5.65	108.66	106.40
36	5	3351	U	N1-C2-O2	5.65	126.76	122.80
1	2	1456	C	N1-C2-O2	5.65	122.29	118.90
36	1	3095	U	C6-N1-C2	-5.65	117.61	121.00
36	1	2373	A	O5'-P-OP1	-5.64	100.62	105.70
36	5	2938	G	O5'-P-OP1	-5.63	100.64	105.70
1	2	720	G	P-O3'-C3'	5.62	126.45	119.70
36	5	3197	G	N3-C4-N9	-5.62	122.63	126.00
36	1	3277	U	N3-C2-O2	-5.61	118.27	122.20
36	5	1222	G	P-O3'-C3'	5.61	126.43	119.70
36	5	2169	G	N3-C4-C5	-5.61	125.80	128.60
36	5	806	A	O5'-P-OP1	-5.60	100.66	105.70
36	1	1269	U	C2-N1-C1'	5.60	124.42	117.70
36	1	1604	G	C4-N9-C1'	5.59	133.77	126.50
44	17	229	PHE	CB-CG-CD2	-5.59	116.89	120.80
36	5	3309	G	N3-C4-C5	-5.59	125.81	128.60
1	2	639	U	N3-C2-O2	-5.57	118.30	122.20
36	5	2938	G	C6-C5-N7	5.57	133.74	130.40
1	2	720	G	OP1-P-O3'	5.57	117.45	105.20
1	2	1600	A	O4'-C1'-N9	5.56	112.65	108.20
78	q2	17	CYS	CA-CB-SG	5.55	124.00	114.00
36	5	942	U	N3-C4-O4	5.55	123.28	119.40
36	1	1151	U	C6-N1-C2	-5.53	117.68	121.00
36	5	1028	U	C6-N1-C1'	-5.53	113.46	121.20
1	6	321	C	N1-C2-O2	5.53	122.22	118.90
36	1	2714	G	N3-C4-N9	-5.52	122.69	126.00
36	5	1152	G	N3-C2-N2	-5.52	116.03	119.90
36	1	439	C	C6-N1-C1'	-5.52	114.18	120.80
36	1	2935	U	N1-C2-O2	5.52	126.66	122.80
36	1	2833	A	C8-N9-C4	5.51	108.00	105.80
36	1	2867	C	N3-C4-C5	5.50	124.10	121.90
36	1	2971	A	P-O3'-C3'	5.50	126.30	119.70
36	1	1434	G	C4-C5-N7	5.49	113.00	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	1	644	G	C5-C6-N1	-5.48	108.76	111.50
1	2	1052	U	N1-C2-O2	5.47	126.63	122.80
48	m1	12	LEU	CA-CB-CG	5.47	127.87	115.30
36	1	2726	C	N3-C2-O2	-5.46	118.08	121.90
36	1	1484	U	P-O3'-C3'	5.46	126.26	119.70
1	2	1052	U	N3-C2-O2	-5.46	118.38	122.20
36	5	942	U	C6-N1-C2	-5.45	117.73	121.00
36	5	1495	U	C5-C6-N1	5.45	125.43	122.70
36	5	421	G	C6-C5-N7	-5.44	127.14	130.40
36	1	2726	C	C6-N1-C2	-5.43	118.13	120.30
36	5	2887	A	N1-C6-N6	5.43	121.86	118.60
36	5	1483	G	O4'-C1'-N9	5.43	112.54	108.20
36	5	1604	G	C4-N9-C1'	5.43	133.56	126.50
24	D2	65	LEU	CA-CB-CG	5.43	127.78	115.30
36	1	1113	G	C5-C6-O6	-5.42	125.35	128.60
1	2	577	G	N1-C6-O6	5.42	123.15	119.90
36	1	2983	C	C2-N1-C1'	5.42	124.76	118.80
36	1	2873	U	N3-C2-O2	-5.41	118.41	122.20
36	1	344	A	N1-C6-N6	-5.41	115.35	118.60
36	1	1419	A	O5'-P-OP1	5.41	117.19	110.70
36	1	1115	G	N9-C4-C5	-5.40	103.24	105.40
36	5	2724	U	C5-C4-O4	5.40	129.14	125.90
1	6	965	U	N1-C2-O2	5.40	126.58	122.80
36	1	2836	C	N1-C2-O2	5.39	122.14	118.90
1	2	1241	G	O4'-C1'-N9	5.39	112.51	108.20
36	1	2298	U	O4'-C1'-N1	5.39	112.51	108.20
36	1	3217	C	C6-N1-C2	-5.39	118.15	120.30
41	14	340	GLY	N-CA-C	-5.38	99.64	113.10
36	1	1820	U	OP2-P-O3'	5.38	117.04	105.20
36	1	2355	G	C5-C6-O6	-5.38	125.37	128.60
31	D9	36	LEU	CA-CB-CG	5.38	127.67	115.30
1	6	144	U	N3-C2-O2	-5.38	118.44	122.20
36	5	419	G	C5-C6-O6	5.37	131.82	128.60
36	1	24	G	O5'-P-OP2	-5.37	100.87	105.70
36	1	503	C	C6-N1-C2	5.36	122.44	120.30
1	2	1456	C	N3-C2-O2	-5.36	118.15	121.90
36	1	65	A	C8-N9-C4	-5.36	103.66	105.80
1	2	992	A	O4'-C1'-N9	5.35	112.48	108.20
1	6	1274	C	N3-C2-O2	-5.35	118.15	121.90
1	6	965	U	C2-N1-C1'	5.34	124.11	117.70
36	5	1208	U	C5-C4-O4	5.34	129.10	125.90
36	5	1368	U	C5-C6-N1	5.33	125.37	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	581	U	C5-C4-O4	-5.33	122.70	125.90
1	6	453	U	C5-C6-N1	5.33	125.36	122.70
36	5	406	G	N3-C4-N9	-5.33	122.81	126.00
36	1	2659	G	C8-N9-C4	5.32	108.53	106.40
36	5	2955	U	OP2-P-O3'	5.32	116.90	105.20
36	1	2112	U	P-O3'-C3'	5.31	126.08	119.70
36	1	2541	U	P-O3'-C3'	5.31	126.07	119.70
36	5	2531	C	C2-N1-C1'	5.31	124.64	118.80
1	2	1761	U	N3-C2-O2	-5.30	118.49	122.20
36	1	1879	A	O4'-C1'-N9	5.30	112.44	108.20
36	5	2273	G	C8-N9-C4	5.30	108.52	106.40
36	5	2964	G	N1-C6-O6	-5.30	116.72	119.90
36	5	2719	U	C2-N1-C1'	-5.28	111.36	117.70
36	1	2297	U	P-O3'-C3'	5.28	126.04	119.70
36	5	3058	U	C2-N1-C1'	5.28	124.03	117.70
36	5	2371	G	C4-N9-C1'	-5.28	119.64	126.50
36	5	3362	A	O4'-C1'-N9	5.28	112.42	108.20
1	2	831	U	N1-C2-O2	5.27	126.49	122.80
36	5	2572	C	C6-N1-C1'	-5.27	114.48	120.80
15	C3	22	ALA	C-N-CD	-5.27	109.02	120.60
1	6	98	U	N3-C2-O2	-5.26	118.52	122.20
36	5	1189	C	N1-C2-O2	-5.26	115.74	118.90
36	5	2827	U	C2-N1-C1'	5.26	124.01	117.70
1	6	275	C	N1-C2-O2	5.26	122.06	118.90
1	6	426	G	C4-N9-C1'	5.26	133.33	126.50
1	2	1657	U	P-O3'-C3'	5.25	126.00	119.70
1	2	934	C	N3-C2-O2	-5.24	118.23	121.90
36	5	1496	C	C2-N1-C1'	5.23	124.55	118.80
36	5	817	A	O5'-P-OP1	-5.22	101.00	105.70
1	2	1000	C	C2-N1-C1'	5.22	124.55	118.80
1	6	558	U	P-O3'-C3'	5.22	125.96	119.70
1	2	639	U	N1-C2-O2	5.22	126.45	122.80
1	6	158	U	P-O3'-C3'	5.22	125.96	119.70
36	5	2205	U	O4'-C1'-N1	5.22	112.37	108.20
36	5	2957	G	OP1-P-OP2	5.21	127.42	119.60
57	N1	89	LEU	CA-CB-CG	5.20	127.25	115.30
36	1	65	A	N7-C8-N9	5.19	116.40	113.80
1	2	1246	C	C2-N1-C1'	5.19	124.51	118.80
1	2	25	C	P-O3'-C3'	5.19	125.93	119.70
36	5	2600	C	C6-N1-C2	-5.19	118.22	120.30
36	5	638	C	C6-N1-C2	-5.19	118.23	120.30
36	1	1269	U	N3-C2-O2	-5.18	118.57	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	5	3145	C	C2-N1-C1'	5.18	124.50	118.80
1	6	453	U	C6-N1-C1'	-5.18	113.95	121.20
36	1	1115	G	C5-C6-O6	-5.18	125.49	128.60
36	5	283	G	C4-C5-N7	5.18	112.87	110.80
36	5	3245	A	C5-N7-C8	-5.18	101.31	103.90
36	5	3140	G	C6-C5-N7	-5.17	127.30	130.40
36	1	1365	G	N3-C4-C5	-5.17	126.02	128.60
36	1	1556	C	C2-N1-C1'	5.17	124.49	118.80
36	5	1166	G	C5-N7-C8	5.17	106.88	104.30
1	2	1389	C	C6-N1-C2	-5.16	118.23	120.30
36	5	663	C	C6-N1-C2	-5.16	118.23	120.30
36	5	2945	G	C5-C6-O6	-5.16	125.50	128.60
1	6	965	U	N3-C2-O2	-5.16	118.59	122.20
47	m0	167	LEU	CA-CB-CG	5.16	127.16	115.30
36	5	3351	U	C5-C6-N1	5.16	125.28	122.70
9	S7	31	SER	C-N-CD	5.15	139.22	128.40
62	N6	126	LEU	CA-CB-CG	5.15	127.15	115.30
36	1	2554	A	P-O3'-C3'	5.15	125.88	119.70
35	SM	134	ASP	CB-CG-OD2	5.15	122.94	118.30
36	1	2690	G	C4-N9-C1'	5.15	133.19	126.50
36	5	661	G	N3-C4-N9	5.15	129.09	126.00
36	1	917	A	N1-C6-N6	-5.15	115.51	118.60
36	5	2772	C	P-O3'-C3'	5.15	125.88	119.70
1	6	1458	G	C4-N9-C1'	5.14	133.19	126.50
36	5	1017	C	N1-C2-O2	5.14	121.99	118.90
36	5	2234	G	C5-C6-N1	-5.14	108.93	111.50
36	1	1349	G	O4'-C1'-N9	5.14	112.31	108.20
36	5	2347	U	N3-C4-O4	-5.14	115.80	119.40
36	5	3317	U	P-O3'-C3'	5.14	125.87	119.70
36	1	1903	U	N3-C4-O4	5.14	123.00	119.40
36	1	3275	U	C5-C6-N1	5.14	125.27	122.70
36	5	1208	U	N3-C4-O4	-5.14	115.80	119.40
36	1	2816	G	O5'-P-OP2	-5.13	101.08	105.70
36	5	1379	G	N3-C2-N2	5.13	123.49	119.90
1	2	1634	C	C2-N1-C1'	5.12	124.43	118.80
36	1	1434	G	C5-N7-C8	-5.11	101.75	104.30
36	5	2945	G	O5'-P-OP2	-5.11	101.10	105.70
1	6	1573	A	OP2-P-O3'	5.10	116.42	105.20
36	5	922	U	C5-C6-N1	-5.10	120.15	122.70
36	5	953	G	C5-C6-O6	-5.10	125.54	128.60
1	2	934	C	C6-N1-C1'	-5.10	114.69	120.80
1	6	1747	G	C8-N9-C4	5.10	108.44	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	6	163	G	N3-C4-C5	5.09	131.15	128.60
1	6	1600	A	OP1-P-O3'	5.09	116.41	105.20
36	1	2101	C	P-O3'-C3'	5.09	125.81	119.70
36	1	1716	U	P-O3'-C3'	5.09	125.81	119.70
36	1	880	G	C4-N9-C1'	-5.09	119.88	126.50
1	6	1389	C	N1-C2-O2	5.09	121.95	118.90
36	1	2617	U	C5-C6-N1	-5.08	120.16	122.70
36	5	968	G	O5'-P-OP1	-5.08	101.13	105.70
36	5	1352	A	P-O3'-C3'	5.08	125.80	119.70
36	5	1604	G	N3-C4-C5	-5.08	126.06	128.60
36	5	2169	G	N1-C6-O6	-5.08	116.86	119.90
1	6	163	G	N3-C2-N2	-5.07	116.35	119.90
1	6	1000	C	C6-N1-C1'	-5.07	114.71	120.80
1	6	217	A	P-O3'-C3'	5.07	125.78	119.70
1	2	577	G	C5-C6-O6	-5.07	125.56	128.60
36	5	1604	G	N3-C4-N9	5.06	129.04	126.00
36	5	27	C	C6-N1-C2	-5.05	118.28	120.30
4	s2	113	LEU	CA-CB-CG	5.04	126.90	115.30
36	1	915	A	C8-N9-C4	-5.04	103.78	105.80
36	1	343	U	O5'-P-OP2	-5.04	101.17	105.70
36	5	1433	A	N9-C4-C5	5.04	107.81	105.80
1	6	144	U	N1-C2-O2	5.03	126.32	122.80
36	5	2156	C	C6-N1-C2	5.03	122.31	120.30
36	1	656	A	C6-C5-N7	-5.02	128.78	132.30
36	5	1116	G	OP2-P-O3'	5.02	116.25	105.20
36	1	2617	U	N3-C2-O2	-5.02	118.69	122.20
36	5	2776	C	N1-C2-O2	5.02	121.91	118.90
47	m0	87	LEU	CA-CB-CG	5.02	126.84	115.30
38	4	125	U	C5-C6-N1	5.02	125.21	122.70
18	C6	40	GLU	C-N-CA	5.01	143.06	122.00
36	5	963	G	O5'-P-OP2	-5.01	101.19	105.70
1	2	767	U	N1-C2-O2	5.01	126.31	122.80
36	1	1269	U	N1-C2-O2	5.01	126.31	122.80
36	1	2971	A	C6-C5-N7	-5.01	128.80	132.30
1	6	163	G	N9-C4-C5	5.00	107.40	105.40
36	1	2137	U	C2-N1-C1'	5.00	123.70	117.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
19	C7	85	VAL	Peptide

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Mol	Chain	Res	Type	Group
27	D5	94	LYS	Peptide
44	L7	157	ASN	Peptide
52	M6	110	PRO	Peptide
65	N9	20	GLY	Peptide
18	c6	40	GLU	Peptide
83	f	108	ASP	Peptide
39	l2	237	LEU	Peptide
64	n8	46	ASP	Peptide
65	n9	23	LYS	Peptide

## 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	2	37970	0	19106	823	0
1	6	38260	0	19252	792	0
2	S0	1577	0	1567	142	0
2	s0	1583	0	1578	0	0
3	S1	1709	0	1784	128	0
3	s1	1722	0	1793	0	0
4	S2	1635	0	1723	138	0
4	s2	1635	0	1723	0	0
5	S3	1734	0	1817	116	0
5	s3	1734	0	1817	0	0
6	S4	2068	0	2154	182	0
6	s4	2068	0	2154	0	0
7	S5	1609	0	1675	118	0
7	s5	1609	0	1675	0	0
8	S6	1799	0	1879	109	0
8	s6	1755	0	1846	0	0
9	S7	1481	0	1572	114	0
9	s7	1491	0	1578	0	0
10	S8	1489	0	1525	114	0
10	s8	1489	0	1525	0	0
11	S9	1494	0	1573	151	0
11	s9	1494	0	1573	0	0
12	C0	773	0	714	57	0
12	c0	762	0	689	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	C1	1214	0	1237	85	0
13	c1	1168	0	1226	0	0
14	C2	892	0	872	64	0
14	c2	892	0	869	0	0
15	C3	1192	0	1255	90	0
15	c3	1192	0	1255	0	0
16	C4	891	0	883	78	0
16	c4	949	0	985	0	0
17	C5	977	0	1004	82	0
17	c5	1039	0	1041	0	0
18	C6	1105	0	1166	85	0
18	c6	1111	0	1171	0	0
19	C7	926	0	930	85	0
19	c7	906	0	909	0	0
20	C8	1192	0	1222	100	0
20	c8	1192	0	1222	0	0
21	C9	1112	0	1124	86	0
21	c9	1112	0	1124	0	0
22	D0	855	0	917	71	0
22	d0	882	0	939	0	0
23	D1	684	0	672	75	0
23	d1	684	0	672	0	0
24	D2	1021	0	1060	97	0
24	d2	1021	0	1060	0	0
25	D3	1121	0	1196	90	0
25	d3	1121	0	1196	0	0
26	D4	1073	0	1132	91	0
26	d4	1073	0	1132	0	0
27	D5	563	0	603	52	0
27	d5	558	0	598	0	0
28	D6	769	0	814	80	0
28	d6	769	0	814	0	0
29	D7	610	0	630	41	0
29	d7	610	0	631	0	0
30	D8	497	0	535	43	0
30	d8	497	0	535	0	0
31	D9	442	0	428	32	0
31	d9	442	0	428	0	0
32	E0	475	0	525	29	0
32	e0	491	0	542	0	0
33	E1	566	0	602	49	0
33	e1	608	0	657	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
34	SR	2437	0	2386	150	0
34	sR	2442	0	2392	0	1
35	SM	1105	0	959	67	0
35	sM	680	0	542	0	0
36	1	67355	0	33848	1240	1
36	5	67780	0	34065	1205	0
37	3	2579	0	1304	61	0
37	7	2579	0	1304	46	0
38	4	3353	0	1695	59	0
38	8	3353	0	1695	70	0
39	L2	1914	0	1981	172	0
39	l2	1912	0	1976	0	0
40	L3	3075	0	3142	240	0
40	l3	3075	0	3142	0	0
41	L4	2748	0	2859	219	0
41	l4	2748	0	2859	0	0
42	L5	2375	0	2325	188	0
42	l5	2359	0	2311	0	0
43	L6	1239	0	1326	79	0
43	l6	1248	0	1339	0	0
44	L7	1784	0	1862	131	0
44	l7	1791	0	1869	0	0
45	L8	1804	0	1875	114	0
45	l8	1764	0	1821	0	0
46	L9	1518	0	1587	120	0
46	l9	1518	0	1587	0	0
47	M0	1705	0	1736	146	0
47	m0	1722	0	1755	0	0
48	M1	1353	0	1383	81	0
48	m1	1353	0	1383	0	0
49	M3	1543	0	1608	137	0
49	m3	1548	0	1613	0	0
50	M4	1053	0	1149	82	0
50	m4	1059	0	1154	0	0
51	M5	1720	0	1779	139	0
51	m5	1720	0	1779	0	0
52	M6	1555	0	1659	106	0
52	m6	1555	0	1659	0	0
53	M7	1420	0	1437	100	0
53	m7	1227	0	1236	0	0
54	M8	1441	0	1543	108	0
54	m8	1441	0	1543	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
55	M9	1521	0	1617	112	0
55	m9	1521	0	1617	0	0
56	N0	1445	0	1487	98	0
56	n0	1445	0	1487	0	0
57	N1	1276	0	1323	99	0
57	n1	1276	0	1323	0	0
58	N2	796	0	812	38	0
58	n2	778	0	791	0	0
59	N3	1003	0	1048	72	0
59	n3	1003	0	1048	0	0
60	N4	699	0	640	43	0
60	n4	1038	0	1071	0	0
61	N5	964	0	1025	71	0
61	n5	959	0	1023	0	0
62	N6	993	0	1081	79	0
62	n6	993	0	1081	0	0
63	N7	1092	0	1155	93	0
63	n7	1092	0	1155	0	0
64	N8	1173	0	1215	108	0
64	n8	1173	0	1215	0	0
65	N9	462	0	491	34	0
65	n9	462	0	491	0	0
66	O0	743	0	797	51	0
66	o0	767	0	816	0	0
67	O1	876	0	912	60	0
67	o1	883	0	918	0	0
68	O2	1020	0	1090	76	0
68	o2	1020	0	1090	0	0
69	O3	850	0	880	64	0
69	o3	850	0	880	0	0
70	O4	880	0	945	69	0
70	o4	880	0	945	0	0
71	O5	969	0	1078	92	0
71	o5	965	0	1067	0	0
72	O6	771	0	849	69	0
72	o6	770	0	846	0	0
73	O7	681	0	683	53	0
73	o7	681	0	683	0	0
74	O8	612	0	682	42	0
74	o8	608	0	671	0	0
75	O9	436	0	475	33	0
75	o9	436	0	475	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
76	Q0	417	0	455	23	0
76	q0	417	0	455	0	0
77	Q1	233	0	284	22	0
77	q1	233	0	284	0	0
78	Q2	847	0	914	50	0
78	q2	847	0	914	0	0
79	Q3	694	0	734	56	0
79	q3	694	0	734	0	0
80	m2	750	0	185	0	0
81	p0	1077	0	1012	0	0
82	p1	235	0	53	0	0
82	p2	230	0	52	0	0
83	f	1116	0	1108	0	0
84	B	76	0	49	3	0
84	C	73	0	49	4	0
85	D6	1	0	0	0	0
85	D7	1	0	0	0	0
85	D9	1	0	0	0	0
85	E1	1	0	0	0	0
85	O7	1	0	0	0	0
85	Q0	1	0	0	0	0
85	Q2	1	0	0	0	0
85	Q3	1	0	0	0	0
85	d6	1	0	0	0	0
85	d7	1	0	0	0	0
85	d9	1	0	0	0	0
85	e1	1	0	0	0	0
85	o7	1	0	0	0	0
85	q0	1	0	0	0	0
85	q2	1	0	0	0	0
85	q3	1	0	0	0	0
86	1	23	0	18	2	0
86	B	23	0	18	3	0
87	1	3	0	0	0	0
87	5	5	0	0	0	0
87	B	1	0	0	0	0
87	C	1	0	0	0	0
87	f	2	0	0	0	0
88	1	14	0	0	0	0
88	5	7	0	0	0	0
89	5	2	0	0	0	0
89	f	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	404292	0	298868	9280	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (9280) 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)
1:6:104:A:H61	1:6:308:C:H5'	1.29	0.96
1:2:1456:C:H5''	1:2:1457:C:H5''	1.49	0.94
34:SR:82:SER:HG	34:SR:92:TRP:HE1	2.15	0.94
36:5:438:A:N1	36:5:621:A:N6	2.17	0.92
36:5:3274:A:H3'	36:5:3275:U:H5''	1.51	0.92
36:5:2620:G:H1	84:B:74:CH:HN3	1.18	0.92
36:1:768:C:OP1	49:M3:186:ARG:NH2	2.02	0.91
5:S3:94:ARG:NH2	35:SM:134:ASP:OD1	2.01	0.91
36:1:2392:C:O2'	40:L3:266:ARG:NH2	2.02	0.91
40:L3:2:SER:N	36:5:2940:A:N7	237.98	0.91
36:5:420:G:N2	36:5:2385:G:OP2	2.04	0.90
62:N6:45:ILE:HD11	62:N6:122:LYS:HB2	1.94	0.90
64:N8:21:ARG:NH2	36:5:640:U:OP1	182.14	0.90
36:1:2656:A:H4'	78:Q2:98:LYS:HD2	1.50	0.90
78:Q2:45:ARG:NH2	36:5:283:G:OP2	147.30	0.89
49:M3:63:VAL:HG22	36:5:72:C:H5'	113.82	0.89
50:M4:55:ARG:NH2	50:M4:76:ALA:O	2.28	0.89
40:L3:232:ARG:NH1	40:L3:269:GLN:O	2.06	0.88
1:6:1636:C:H4'	1:6:1637:C:H5'	1.56	0.88
62:N6:5:SER:HB2	62:N6:8:VAL:HG22	7.22	0.87
41:L4:112:LYS:HD2	36:5:790:U:H5'	121.05	0.87
11:S9:96:VAL:HA	11:S9:99:LEU:HD23	1.54	0.87
51:M5:188:ARG:NH2	36:5:31:C:OP2	121.33	0.87
2:S0:162:CYS:SG	2:S0:163:ASN:N	2.47	0.87
39:L2:21:ARG:HD3	36:5:824:C:H5''	170.23	0.86
36:1:640:U:OP1	64:N8:21:ARG:NH2	2.09	0.86
1:2:992:A:H2	1:2:1012:U:H3	1.23	0.86
36:1:860:G:OP1	79:Q3:17:ARG:NH1	2.09	0.85
39:L2:153:GLY:HA3	39:L2:251:LYS:HD2	7.65	0.85
47:M0:42:THR:HG23	47:M0:45:GLU:HG3	4.60	0.85
52:M6:108:ILE:HD12	52:M6:113:ASP:HA	4.75	0.85
6:S4:9:LEU:HB2	6:S4:30:ARG:HB2	2.30	0.85
41:L4:329:PRO:O	41:L4:331:ALA:N	3.21	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:14:LYS:HA	51:M5:19:LEU:HD23	1.59	0.85
10:S8:84:HIS:HE2	10:S8:97:THR:HG1	1.55	0.85
57:N1:51:GLY:HA3	57:N1:92:ARG:HG3	2.07	0.84
36:5:3299:A:H61	36:5:3315:G:H1	1.24	0.84
36:1:687:U:OP2	49:M3:36:ARG:NH2	2.09	0.84
19:C7:66:VAL:HB	19:C7:69:ILE:HD11	1.60	0.84
20:C8:27:LYS:HA	20:C8:57:ARG:HA	2.35	0.84
8:S6:98:ARG:NH2	8:S6:101:ILE:O	2.18	0.84
40:L3:41:VAL:HA	40:L3:185:GLY:HA3	1.71	0.84
48:M1:94:ARG:O	48:M1:96:PHE:N	2.77	0.84
41:L4:300:ARG:O	54:M8:39:ARG:NH1	2.11	0.84
18:C6:9:THR:HG21	18:C6:88:GLY:HA2	1.59	0.83
10:S8:34:ALA:HB2	10:S8:56:ARG:HD3	2.16	0.83
36:1:3042:U:OP2	36:1:3092:C:N4	2.11	0.83
18:C6:46:PHE:HA	18:C6:49:TYR:HB2	1.59	0.83
3:S1:51:SER:HB3	3:S1:57:ALA:H	3.96	0.83
1:2:630:A:N6	1:2:969:C:O2	2.12	0.83
39:L2:128:ARG:NH1	36:5:2177:G:OP2	198.84	0.83
17:C5:122:THR:HG22	1:6:1558:U:H3	367.07	0.82
3:S1:144:ARG:HB3	3:S1:208:GLN:HG2	4.20	0.82
6:S4:206:ASP:HB2	6:S4:222:LEU:HD12	2.15	0.82
57:N1:87:LYS:NZ	36:5:2728:G:N7	212.25	0.82
51:M5:143:ARG:HH21	71:O5:92:LEU:HD23	1.42	0.82
36:1:3182:G:OP1	52:M6:160:ARG:NH2	2.11	0.82
50:M4:108:ARG:NH2	52:M6:196:ALA:O	2.12	0.82
7:S5:112:ARG:HH11	18:C6:43:ILE:HG23	5.27	0.82
47:M0:43:VAL:HG21	47:M0:197:VAL:HB	2.58	0.82
36:1:837:A:OP2	79:Q3:4:ARG:NH1	2.13	0.82
36:1:2940:A:N7	40:L3:2:SER:N	2.27	0.82
40:L3:169:THR:O	40:L3:171:LEU:N	2.97	0.82
11:S9:81:VAL:HG22	11:S9:86:LEU:HB3	3.08	0.82
20:C8:143:ARG:NH2	1:6:1462:G:N7	338.70	0.82
36:1:2771:U:OP2	36:1:2772:C:N4	2.13	0.81
1:2:1096:C:H4'	1:2:1099:U:H4'	1.62	0.81
68:O2:44:ARG:NH1	36:5:1145:G:OP1	206.95	0.81
63:N7:135:ARG:NH2	36:5:1807:G:OP1	196.98	0.81
44:L7:151:ARG:NH1	44:L7:244:ASN:O	3.54	0.81
17:C5:18:ARG:NH1	20:C8:91:ASP:O	6.01	0.81
36:1:1639:C:OP2	70:O4:74:ARG:NH2	2.13	0.81
1:2:1537:C:O2'	1:2:1540:G:O6	1.98	0.81
36:5:1235:U:H4'	36:5:1236:G:H5'	1.62	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1034:C:HO2'	24:D2:2:THR:N	1.80	0.80
36:5:2836:C:H5	36:5:2852:C:H42	1.27	0.80
1:6:647:G:H1	1:6:687:G:H22	1.29	0.80
17:C5:123:TYR:OH	20:C8:126:ARG:NH1	2.14	0.80
23:D1:74:GLN:NE2	23:D1:81:ASN:O	2.14	0.80
36:5:155:G:H5''	36:5:156:G:C8	2.17	0.80
14:C2:103:LEU:HG	14:C2:116:VAL:HG22	1.63	0.80
4:S2:77:GLN:NE2	4:S2:105:GLY:O	2.14	0.80
8:S6:14:LYS:HD2	8:S6:123:GLY:HA3	2.29	0.80
17:C5:43:ARG:NH2	1:6:1552:U:OP2	402.72	0.80
10:S8:176:SER:HB3	1:6:208:U:H4'	286.17	0.80
36:1:2619:G:H1	84:C:75:CH:HN3	1.29	0.80
64:N8:67:HIS:NE2	36:5:71:A:OP2	119.15	0.80
18:C6:110:THR:HA	18:C6:113:ASP:HB2	3.02	0.80
39:L2:29:LEU:HA	39:L2:76:PHE:HE1	1.45	0.80
62:N6:56:VAL:HG11	62:N6:104:LEU:HD13	2.43	0.80
1:2:1542:G:N2	1:2:1569:A:OP2	2.15	0.80
39:L2:152:SER:OG	36:5:2157:G:N7	216.53	0.80
57:N1:39:ILE:HD12	57:N1:102:ARG:HG2	1.64	0.80
15:C3:65:VAL:HG23	15:C3:66:ILE:HG23	5.87	0.80
42:L5:40:HIS:HD2	42:L5:42:ALA:H	1.30	0.80
6:S4:52:LEU:O	6:S4:54:TYR:N	2.74	0.80
36:5:2812:C:H2'	36:5:2813:A:H8	1.46	0.80
1:6:219:A:N6	1:6:843:U:O2	2.14	0.80
52:M6:62:THR:H	52:M6:69:GLY:HA3	1.75	0.79
68:O2:4:LEU:HD12	68:O2:5:PRO:HD2	1.64	0.79
3:S1:134:VAL:HB	3:S1:219:LYS:HB2	1.64	0.79
8:S6:26:VAL:HG21	8:S6:40:ALA:HB1	1.63	0.79
40:L3:347:SER:OG	40:L3:348:ARG:N	2.13	0.79
71:O5:10:ARG:NH1	71:O5:60:GLU:OE1	4.56	0.79
35:SM:53:ARG:HE	35:SM:54:PRO:HD2	1.48	0.79
38:4:95:G:OP2	73:O7:72:ARG:NH1	2.15	0.79
21:C9:52:GLY:O	21:C9:54:PHE:N	2.15	0.79
70:O4:9:ARG:HH21	70:O4:34:HIS:HB2	2.64	0.79
35:SM:48:ARG:HH21	36:1:1017:C:H5'	1.47	0.79
36:1:977:C:OP1	54:M8:141:ARG:NH2	2.15	0.79
1:2:1083:G:HO2'	1:2:1094:G:HO2'	1.26	0.79
41:L4:16:THR:HG23	41:L4:18:ASN:H	2.21	0.79
34:SR:24:ALA:HB3	34:SR:34:LEU:HB3	2.19	0.79
36:1:412:G:OP1	53:M7:62:ARG:NH1	2.15	0.79
41:L4:122:THR:HG22	41:L4:235:LEU:HB2	1.88	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:3194:C:O2	36:5:3197:G:N2	2.16	0.78
36:1:994:G:H3'	57:N1:13:TYR:HD2	1.48	0.78
1:6:680:U:O2	1:6:682:C:N4	2.16	0.78
71:O5:78:LYS:HA	71:O5:81:ARG:HD2	1.84	0.78
25:D3:92:CYS:HA	25:D3:95:PHE:HD2	1.49	0.78
66:O0:16:LEU:HB2	66:O0:98:SER:HB2	2.13	0.78
24:D2:103:ILE:HA	24:D2:112:ASP:HA	2.05	0.78
53:M7:62:ARG:NH1	36:5:412:G:OP1	159.33	0.78
63:N7:83:THR:HG23	63:N7:85:TYR:H	1.48	0.78
1:2:1793:G:N2	28:D6:76:SER:OG	2.16	0.78
36:5:3241:G:H2'	36:5:3245:A:C8	2.19	0.78
65:N9:14:ARG:HH12	65:N9:18:ARG:HH11	1.31	0.78
36:5:2569:A:H4'	36:5:2570:U:H5'	1.66	0.78
53:M7:111:LYS:HE2	53:M7:152:GLU:HB3	4.56	0.78
1:2:1188:G:O2'	1:2:1430:U:OP1	2.02	0.78
1:6:709:C:O2	1:6:730:G:N2	2.17	0.78
36:1:398:A:H5''	53:M7:3:ARG:HD2	1.66	0.77
1:2:1595:U:H3	1:2:1600:A:H2	1.31	0.77
56:N0:115:ARG:NH1	36:5:1295:G:O2'	294.48	0.77
52:M6:179:ALA:HA	52:M6:182:ASN:HB3	5.18	0.77
70:O4:80:ARG:HG3	70:O4:88:ARG:HH21	3.09	0.77
36:1:542:G:H1	36:1:549:U:H3	1.29	0.77
1:2:930:A:OP1	28:D6:32:LYS:NZ	2.17	0.77
41:L4:269:SER:O	41:L4:271:LYS:N	2.15	0.77
1:6:119:A:N6	1:6:298:C:O2	2.17	0.77
36:1:1553:U:H4'	36:1:1554:U:H5'	1.64	0.77
46:L9:49:ASN:O	46:L9:51:GLN:N	2.18	0.77
53:M7:50:GLN:OE1	53:M7:56:ARG:NH2	2.18	0.77
36:1:709:A:OP1	54:M8:179:ARG:NH2	2.17	0.77
75:O9:23:LEU:O	75:O9:25:GLN:NE2	2.67	0.77
1:2:1067:C:H5''	3:S1:150:VAL:HG23	1.67	0.77
64:N8:19:LYS:HD2	64:N8:25:HIS:HD2	5.16	0.77
64:N8:6:THR:HG23	64:N8:8:THR:HG23	1.66	0.77
36:5:2534:G:H1	36:5:2545:C:H42	1.32	0.77
40:L3:260:VAL:HG11	40:L3:266:ARG:NH1	2.00	0.77
41:L4:292:SER:OG	41:L4:293:SER:N	2.18	0.77
26:D4:116:LYS:NZ	1:6:57:G:OP2	338.44	0.77
42:L5:22:ARG:NH2	42:L5:28:THR:OG1	2.16	0.77
52:M6:12:LYS:O	56:N0:167:ARG:NH2	2.18	0.77
36:1:3245:A:H5'	36:1:3246:G:H5''	1.66	0.77
64:N8:58:MET:SD	36:5:2786:G:N2	156.07	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D6:23:CYS:SG	28:D6:74:CYS:HB3	2.24	0.76
4:S2:101:VAL:HG22	4:S2:115:ILE:HG22	3.59	0.76
36:1:824:C:H5''	39:L2:21:ARG:HD3	1.66	0.76
46:L9:28:VAL:HG13	46:L9:33:THR:HB	1.66	0.76
1:2:1267:G:HO2'	1:2:1448:G:HO2'	1.26	0.76
13:C1:99:ARG:NH1	25:D3:7:ARG:O	2.31	0.76
36:1:1940:G:H21	36:1:3362:A:H8	1.32	0.76
36:1:655:C:H2'	36:1:656:A:C8	2.19	0.76
13:C1:14:GLN:HB3	13:C1:54:ILE:HG13	3.86	0.76
49:M3:28:GLN:HB3	51:M5:201:ARG:HH11	2.40	0.76
64:N8:47:LYS:O	64:N8:49:HIS:N	2.45	0.76
56:N0:4:PHE:HD2	56:N0:104:GLU:HG3	2.71	0.76
36:1:939:U:OP2	64:N8:26:ARG:NH2	2.19	0.76
41:L4:219:LEU:O	41:L4:221:ASN:N	2.19	0.76
47:M0:174:THR:HG23	47:M0:176:LEU:H	1.49	0.76
36:1:3275:U:H5'	69:O3:68:TRP:HZ2	1.49	0.76
1:2:1559:A:H5''	20:C8:135:GLY:HA3	1.68	0.76
39:L2:181:LYS:HE2	39:L2:184:ARG:HE	1.49	0.76
36:1:2916:U:H1'	59:N3:44:SER:HB3	1.68	0.76
1:2:283:U:H5''	8:S6:188:ARG:HD3	1.68	0.75
48:M1:26:SER:OG	48:M1:27:GLY:N	2.18	0.75
36:5:2211:U:H5	36:5:2234:G:H1	1.34	0.75
46:L9:38:LEU:HD13	46:L9:71:VAL:HG22	1.67	0.75
49:M3:52:ASP:N	49:M3:52:ASP:OD1	2.65	0.75
67:O1:13:THR:HG22	67:O1:72:ARG:HH21	5.84	0.75
36:1:3386:G:OP1	67:O1:10:ARG:NH2	2.19	0.75
2:S0:84:ARG:HH21	2:S0:201:LEU:HA	4.28	0.75
4:S2:56:ILE:HG23	4:S2:61:LEU:HB2	3.73	0.75
67:O1:55:LEU:HB2	67:O1:95:PRO:HD3	2.02	0.75
51:M5:46:ASP:OD1	51:M5:50:ARG:NH2	2.20	0.75
55:M9:104:ARG:NH1	36:5:1949:G:OP1	218.05	0.75
1:6:828:U:H3	1:6:844:A:H61	1.35	0.75
27:D5:71:ILE:HG23	27:D5:73:GLY:H	8.55	0.75
36:1:1318:A:OP1	52:M6:128:ARG:NH1	2.19	0.75
47:M0:41:ALA:O	47:M0:139:ARG:NH2	5.17	0.75
24:D2:15:ASN:ND2	24:D2:72:CYS:SG	5.33	0.75
36:1:1245:A:H3'	36:1:1246:G:H5''	1.68	0.74
36:1:14:U:O4	38:4:136:G:N2	2.19	0.74
1:2:1601:G:OP1	21:C9:86:ARG:NH2	2.20	0.74
40:L3:92:TYR:HB2	40:L3:157:VAL:HG22	1.69	0.74
41:L4:82:THR:O	41:L4:84:ARG:N	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:N0:155:ARG:NH1	36:5:3206:C:O2	310.38	0.74
1:2:1484:G:H21	1:2:1606:C:H1'	1.52	0.74
61:N5:82:LEU:HD11	61:N5:135:ILE:HD11	4.60	0.74
36:1:514:G:N3	41:L4:341:SER:OG	2.20	0.74
37:3:97:A:O4'	44:L7:225:GLN:NE2	2.21	0.74
36:1:2771:U:O2'	36:1:2772:C:O4'	2.04	0.74
36:5:283:G:OP2	36:5:285:A:O2'	2.05	0.74
63:N7:33:SER:OG	63:N7:35:SER:O	5.29	0.74
2:S0:41:ARG:NE	2:S0:42:PRO:O	2.21	0.74
36:1:1555:U:H5	36:1:1559:A:H61	1.35	0.74
12:C0:56:LYS:HB3	12:C0:67:THR:HB	1.70	0.74
42:L5:107:ARG:HH21	42:L5:110:LEU:HD23	1.52	0.74
42:L5:68:THR:HG22	42:L5:70:THR:H	1.51	0.74
1:2:159:U:O2'	8:S6:87:ARG:NH1	2.21	0.74
11:S9:53:ARG:NH2	11:S9:97:LEU:O	2.21	0.74
36:5:1565:G:N1	36:5:1574:C:N3	2.36	0.74
36:1:103:G:OP1	49:M3:70:ARG:NH2	2.21	0.74
1:2:803:A:N3	9:S7:104:ARG:NH2	2.36	0.74
36:5:2568:C:N4	36:5:2574:G:O6	2.18	0.74
23:D1:3:ASN:ND2	23:D1:7:GLN:O	3.53	0.74
25:D3:137:LYS:O	25:D3:139:LYS:N	4.84	0.74
47:M0:169:LYS:HE3	57:N1:159:PHE:HA	4.02	0.74
6:S4:221:ARG:HG3	1:6:753:A:H5'	360.77	0.74
10:S8:84:HIS:NE2	10:S8:97:THR:OG1	2.41	0.74
36:5:252:U:H4'	36:5:253:A:H5'	1.70	0.74
1:6:918:U:H2'	1:6:919:A:H8	1.52	0.74
6:S4:185:GLY:H	6:S4:189:LEU:HB2	1.51	0.74
1:2:246:G:N2	13:C1:38:ALA:O	2.21	0.73
24:D2:80:ASN:OD1	24:D2:124:LYS:NZ	3.15	0.73
41:L4:26:PHE:HD1	41:L4:130:ALA:HB2	1.87	0.73
2:S0:24:LEU:O	2:S0:163:ASN:ND2	2.20	0.73
53:M7:40:GLU:HG2	53:M7:42:THR:HG23	1.70	0.73
56:N0:2:ALA:HB3	56:N0:32:SER:HB3	1.70	0.73
1:2:1773:C:OP2	77:Q1:2:ARG:NH1	2.21	0.73
6:S4:49:ARG:NH2	6:S4:50:ASN:OD1	3.90	0.73
39:L2:224:THR:HG21	36:5:2201:G:H21	222.27	0.73
36:1:116:A:OP1	72:O6:36:ARG:NH1	2.21	0.73
36:5:979:U:H1'	36:5:980:A:C4	2.24	0.73
1:6:224:C:H2'	1:6:225:A:H8	1.53	0.73
40:L3:293:ASN:HB2	40:L3:304:THR:HA	1.68	0.73
47:M0:193:ASP:OD1	47:M0:198:LYS:NZ	2.20	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1603:A:OP1	55:M9:38:ARG:NH1	2.21	0.73
63:N7:46:ILE:HD13	63:N7:68:ILE:HG23	1.78	0.73
5:S3:190:ARG:HH22	5:S3:195:SER:HA	1.54	0.73
1:2:478:A:O2'	11:S9:124:HIS:ND1	2.22	0.73
21:C9:102:ARG:NH2	1:6:1502:G:N7	406.33	0.73
53:M7:25:SER:O	53:M7:29:THR:HG23	1.89	0.73
7:S5:57:SER:O	7:S5:59:VAL:N	2.20	0.73
6:S4:49:ARG:NH1	1:6:448:C:OP2	379.21	0.73
36:5:1352:A:H4'	36:5:1353:U:OP1	1.88	0.73
41:L4:328:ASN:OD1	44:L7:48:ASN:ND2	2.31	0.73
47:M0:156:ARG:NH1	47:M0:163:GLN:O	3.67	0.73
67:O1:64:VAL:HG23	67:O1:65:LYS:HG3	1.69	0.73
10:S8:76:THR:HG22	10:S8:108:PRO:HG2	1.69	0.73
36:1:2213:A:H2'	36:1:2214:A:C8	2.24	0.73
1:2:761:G:OP1	11:S9:54:ARG:NH1	2.22	0.73
38:4:83:C:O2	62:N6:51:ARG:NH2	2.21	0.73
13:C1:91:LEU:HB3	13:C1:100:TYR:HB3	1.69	0.73
22:D0:61:LYS:HB2	22:D0:86:ILE:HB	1.70	0.73
54:M8:158:HIS:H	54:M8:186:VAL:HG12	1.53	0.73
36:1:361:A:O3'	73:O7:45:ARG:NH2	2.22	0.73
11:S9:23:ARG:NH1	11:S9:27:GLU:OE2	3.58	0.73
36:1:3366:G:OP1	60:N4:61:LYS:NZ	2.17	0.73
36:1:718:G:C2	36:1:721:G:H1'	2.24	0.73
18:C6:128:LYS:NZ	18:C6:134:ALA:O	2.22	0.73
1:2:1777:G:O6	77:Q1:8:LYS:NZ	2.22	0.73
4:S2:179:VAL:O	4:S2:198:THR:OG1	2.06	0.73
65:N9:47:LEU:HA	65:N9:50:THR:HG22	1.70	0.73
1:6:1595:U:H3	1:6:1600:A:H2	1.35	0.72
8:S6:163:THR:HG22	8:S6:168:THR:HG22	1.70	0.72
16:C4:131:GLY:O	16:C4:133:ARG:N	2.22	0.72
4:S2:229:LEU:O	23:D1:16:LYS:NZ	2.22	0.72
39:L2:213:GLY:HA3	36:5:2967:A:H5''	204.58	0.72
39:L2:242:ARG:HD3	39:L2:246:LEU:HD12	5.83	0.72
51:M5:144:ARG:NH1	36:5:126:U:OP1	76.92	0.72
10:S8:42:ARG:NH1	1:6:1677:C:OP1	262.96	0.72
11:S9:143:ILE:HD13	1:6:767:U:H5	420.65	0.72
59:N3:3:GLY:HA2	59:N3:40:LYS:HB3	5.90	0.72
8:S6:87:ARG:NH1	1:6:159:U:O2'	321.23	0.72
1:2:652:G:H1	1:2:682:C:H42	1.35	0.72
36:5:1863:G:N1	36:5:1866:C:OP2	2.18	0.72
1:6:52:U:H2'	1:6:53:G:H8	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:472:U:O2'	1:6:769:A:N3	2.22	0.72
62:N6:36:SER:HB3	62:N6:39:LEU:HB2	1.71	0.72
71:O5:30:GLU:OE1	71:O5:34:GLN:NE2	9.10	0.72
11:S9:66:ASP:HB3	11:S9:69:ARG:HB3	2.47	0.72
1:2:901:G:OP2	1:2:901:G:N2	2.20	0.72
51:M5:144:ARG:HE	71:O5:95:PHE:HE2	1.35	0.72
6:S4:251:GLU:HB3	6:S4:255:ARG:HH12	1.54	0.72
14:C2:61:VAL:HG13	14:C2:121:VAL:HG23	3.96	0.72
41:L4:25:VAL:HG22	41:L4:262:TRP:HB2	1.72	0.72
62:N6:23:PRO:HD2	62:N6:26:GLN:HB2	3.19	0.72
3:S1:129:THR:HA	3:S1:177:GLN:HA	1.69	0.72
36:1:1043:C:O3'	47:M0:90:ARG:NH1	2.23	0.72
1:6:884:A:H2'	1:6:885:G:H8	1.54	0.72
40:L3:3:HIS:ND1	40:L3:3:HIS:O	2.21	0.72
36:1:2736:A:OP1	57:N1:92:ARG:NH1	2.22	0.72
70:O4:74:ARG:NH2	36:5:1639:C:OP2	199.49	0.72
8:S6:164:LYS:HB3	8:S6:167:LYS:HB3	1.69	0.72
1:6:1000:C:N4	1:6:1003:A:OP2	2.17	0.72
11:S9:17:ARG:NH1	1:6:4:C:O2'	388.73	0.72
40:L3:166:ILE:HD13	40:L3:173:GLN:HG2	1.71	0.72
44:L7:140:SER:HB2	44:L7:143:THR:H	3.90	0.72
46:L9:112:ILE:HD11	46:L9:134:ILE:HD13	1.71	0.72
64:N8:29:PRO:O	64:N8:31:GLY:N	3.47	0.72
71:O5:85:THR:HG22	71:O5:87:ALA:H	1.53	0.72
4:S2:56:ILE:HA	4:S2:61:LEU:HD12	1.70	0.72
38:4:135:G:H5''	61:N5:49:LYS:HD2	1.72	0.72
42:L5:58:LYS:HD2	42:L5:93:THR:HG21	1.72	0.72
63:N7:127:ASN:O	63:N7:129:TRP:N	2.23	0.72
67:O1:10:ARG:HE	67:O1:108:VAL:HG22	1.52	0.72
36:1:1563:C:O2	36:1:1577:G:N2	2.23	0.72
43:L6:55:LEU:HD11	43:L6:66:SER:HB2	2.32	0.72
36:1:40:A:H5''	64:N8:35:ALA:HB1	1.72	0.72
8:S6:160:ARG:HH12	1:6:68:A:H5'	346.19	0.71
18:C6:50:GLU:OE2	18:C6:82:ARG:NH2	2.23	0.71
64:N8:16:SER:HA	36:5:942:U:N3	170.28	0.71
5:S3:164:VAL:HG13	5:S3:168:ILE:HD11	1.72	0.71
36:1:224:C:O2	62:N6:103:LYS:NZ	2.23	0.71
1:2:885:G:H21	16:C4:123:SER:HB2	1.53	0.71
36:5:3242:G:H5'	36:5:3245:A:H8	1.54	0.71
16:C4:90:ARG:O	16:C4:92:LYS:N	3.07	0.71
25:D3:125:VAL:HG12	25:D3:126:LYS:HG3	4.67	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:L5:177:GLU:O	42:L5:179:ARG:N	2.22	0.71
46:L9:87:LYS:HD3	46:L9:89:LYS:HE3	1.72	0.71
47:M0:5:PRO:O	47:M0:7:ARG:N	2.22	0.71
55:M9:13:SER:OG	55:M9:38:ARG:NH2	2.22	0.71
2:S0:65:ALA:O	2:S0:67:ILE:N	4.74	0.71
36:1:3194:C:O2	36:1:3197:G:N2	2.18	0.71
1:2:1459:C:OP1	20:C8:126:ARG:NH2	2.23	0.71
47:M0:140:THR:HG21	47:M0:144:ASN:HD22	1.56	0.71
49:M3:140:SER:OG	49:M3:143:ALA:N	2.21	0.71
44:L7:90:LYS:NZ	36:5:1158:A:OP2	241.42	0.71
33:E1:91:ILE:HG12	33:E1:92:LYS:HG2	6.52	0.71
43:L6:31:ARG:NH1	69:O3:107:ILE:O	2.23	0.71
75:O9:10:LYS:NZ	36:5:1833:G:OP1	104.57	0.71
36:5:198:A:N3	36:5:218:G:O2'	2.23	0.71
21:C9:31:PRO:HG3	21:C9:103:LYS:HD3	1.71	0.71
24:D2:9:ASP:N	24:D2:9:ASP:OD1	3.53	0.71
48:M1:85:LYS:NZ	48:M1:89:TYR:OH	2.24	0.71
50:M4:37:GLU:OE2	50:M4:74:ARG:NH2	3.88	0.71
35:SM:64:LYS:O	35:SM:66:ALA:N	2.21	0.71
36:1:2353:G:H5''	53:M7:86:LYS:HB2	1.72	0.71
51:M5:31:ARG:NH1	51:M5:124:ASP:OD1	2.25	0.71
1:2:1122:G:N2	1:2:1125:A:OP2	2.23	0.71
33:E1:103:LEU:HD11	1:6:1252:C:H5'	454.29	0.71
16:C4:86:THR:HG21	16:C4:90:ARG:HD2	1.72	0.71
13:C1:101:GLU:OE2	25:D3:16:ARG:NH2	2.99	0.71
44:L7:88:ARG:HG2	44:L7:111:ILE:HA	1.73	0.71
36:5:496:C:O2	36:5:617:G:N2	2.18	0.71
55:M9:62:ARG:NH2	36:5:3068:U:OP2	172.90	0.71
36:1:2203:U:H4'	39:L2:241:ARG:HA	1.71	0.71
36:5:2449:A:H2'	36:5:2450:G:H8	1.55	0.71
28:D6:95:ARG:NH1	1:6:1796:C:O2'	340.95	0.71
1:6:591:A:H2'	1:6:592:A:C8	2.26	0.71
40:L3:250:ALA:HB3	36:5:2880:U:H1'	223.82	0.71
36:1:2828:G:OP1	47:M0:7:ARG:NH1	2.23	0.71
11:S9:108:ARG:HH21	11:S9:145:SER:HB2	1.55	0.71
11:S9:126:ARG:NH1	1:6:475:A:OP2	423.49	0.71
36:1:3343:G:H21	36:1:3362:A:H2	1.38	0.71
1:2:1473:U:O2'	7:S5:103:ASN:ND2	2.23	0.71
1:6:1186:U:O4	1:6:1200:G:N2	2.24	0.71
1:6:158:U:O2'	1:6:160:C:OP2	2.08	0.71
16:C4:97:GLY:O	16:C4:99:GLN:N	4.01	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:3344:A:H2	36:1:3361:G:H21	1.37	0.70
68:O2:19:ARG:HH11	68:O2:28:VAL:HG13	2.93	0.70
6:S4:103:TYR:O	6:S4:182:TYR:OH	2.08	0.70
7:S5:166:ARG:HA	7:S5:169:ASN:HB2	2.73	0.70
5:S3:223:LYS:HD3	34:SR:193:ILE:HD13	3.54	0.70
74:O8:2:ALA:HA	36:5:1747:G:H21	144.64	0.70
16:C4:20:TYR:HB3	16:C4:27:PHE:HB2	1.85	0.70
22:D0:106:ILE:HG23	22:D0:107:THR:HG23	1.73	0.70
40:L3:138:ALA:O	40:L3:140:ASP:N	4.21	0.70
36:1:2880:U:H1'	40:L3:250:ALA:HB3	1.72	0.70
44:L7:41:ARG:NH1	36:5:598:A:OP1	260.51	0.70
4:S2:170:ILE:HB	4:S2:197:TYR:HB2	1.84	0.70
23:D1:60:ARG:HA	23:D1:65:SER:HB2	3.07	0.70
42:L5:40:HIS:HE1	42:L5:42:ALA:HB3	4.32	0.70
46:L9:138:THR:O	46:L9:139:ASN:ND2	2.24	0.70
65:N9:59:LYS:HD3	65:N9:59:LYS:H	1.57	0.70
70:O4:46:ASP:OD1	70:O4:80:ARG:NH1	2.23	0.70
7:S5:145:ASP:OD1	30:D8:45:LYS:NZ	2.22	0.70
7:S5:37:GLN:HB3	18:C6:53:LEU:HD22	1.73	0.70
1:6:514:G:HO2'	1:6:515:A:H8	1.39	0.70
32:E0:13:LYS:HE2	32:E0:17:GLN:HE22	4.06	0.70
36:1:2177:G:OP2	39:L2:128:ARG:NH1	2.25	0.70
42:L5:85:ARG:HH12	42:L5:254:LYS:H	2.09	0.70
36:1:3272:C:OP2	43:L6:78:ARG:NH1	2.23	0.70
58:N2:54:VAL:HG13	58:N2:67:SER:HB2	4.07	0.70
70:O4:87:GLU:OE1	70:O4:91:ARG:NH1	4.23	0.70
34:SR:178:VAL:HB	34:SR:192:PHE:HB2	1.94	0.70
23:D1:39:VAL:HA	23:D1:45:ALA:HA	1.73	0.70
48:M1:6:GLN:O	48:M1:7:ASN:ND2	2.25	0.70
36:1:2818:U:H6	36:1:2818:U:H5'	1.55	0.70
36:1:2898:G:N7	76:Q0:125:LYS:NZ	2.37	0.70
36:5:1141:C:O2'	36:5:1153:A:N3	2.25	0.70
24:D2:119:LYS:HG2	1:6:687:G:H5''	392.25	0.70
45:L8:50:VAL:HG12	61:N5:30:ALA:HA	1.73	0.70
79:Q3:72:SER:OG	79:Q3:80:ARG:NH2	2.23	0.70
3:S1:123:ALA:HB2	3:S1:165:ARG:HG2	1.74	0.70
50:M4:128:ARG:NH2	36:5:3214:U:OP2	280.77	0.70
15:C3:76:LYS:HA	15:C3:81:ALA:HB2	2.57	0.70
28:D6:4:LYS:NZ	1:6:1795:U:OP2	339.97	0.70
44:L7:107:ARG:HH21	44:L7:200:ASN:HA	1.57	0.70
58:N2:50:LEU:HB3	58:N2:54:VAL:HG23	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S2:153:SER:OG	4:S2:154:LEU:N	2.19	0.70
36:1:2986:U:H2'	36:1:2987:A:C8	2.27	0.70
8:S6:164:LYS:NZ	1:6:71:A:OP2	370.75	0.70
1:6:868:G:H1	1:6:960:U:H3	1.37	0.70
17:C5:130:ARG:H	17:C5:130:ARG:HH11	1.40	0.70
27:D5:43:ASP:O	27:D5:46:LYS:N	2.24	0.70
45:L8:140:VAL:HG22	45:L8:166:LEU:HD21	1.74	0.70
51:M5:47:LYS:HE3	51:M5:51:LEU:HD11	1.74	0.70
59:N3:54:LEU:HD21	59:N3:119:GLY:HA3	3.04	0.70
59:N3:66:LYS:HD2	59:N3:68:GLU:HB2	5.36	0.70
11:S9:81:VAL:HG21	11:S9:91:LYS:HE3	1.73	0.70
36:1:1097:G:H5'	57:N1:129:LYS:HE3	1.73	0.70
1:6:151:G:H1	1:6:163:G:H1	1.39	0.70
13:C1:125:VAL:HG12	13:C1:139:VAL:HA	1.96	0.70
13:C1:77:SER:HB3	13:C1:85:VAL:HB	1.72	0.70
42:L5:64:ILE:HD13	42:L5:144:VAL:HG21	4.69	0.70
59:N3:87:ARG:HH22	59:N3:137:VAL:HG22	1.66	0.70
9:S7:154:LEU:HD21	9:S7:183:PHE:HD1	1.56	0.70
11:S9:113:VAL:HG12	11:S9:119:ALA:HB2	4.56	0.70
1:6:1588:G:H1	1:6:1608:U:H3	1.39	0.70
19:C7:67:ARG:NH2	1:6:1398:U:O2'	404.37	0.70
7:S5:64:VAL:HG12	7:S5:65:ARG:HD3	1.74	0.70
34:SR:159:ASN:O	34:SR:161:LYS:N	3.92	0.70
36:5:1952:G:H1	36:5:2094:C:H42	1.40	0.69
10:S8:178:ARG:NH1	1:6:207:U:O2	287.74	0.69
17:C5:43:ARG:NH1	1:6:1553:G:O6	398.91	0.69
49:M3:46:ILE:HG22	49:M3:49:ARG:HB2	2.60	0.69
36:1:436:A:H2'	36:1:437:G:O4'	1.93	0.69
36:5:655:C:H2'	36:5:656:A:H8	1.58	0.69
15:C3:55:ARG:NH1	15:C3:56:ASP:OD2	2.25	0.69
16:C4:50:ALA:O	16:C4:52:ARG:N	2.37	0.69
54:M8:71:LEU:HD13	54:M8:99:THR:HG21	1.74	0.69
10:S8:5:ARG:NH1	10:S8:29:LEU:O	3.08	0.69
36:1:1308:A:C8	36:1:1308:A:OP2	2.45	0.69
36:1:2345:A:OP1	67:O1:24:SER:OG	2.07	0.69
36:1:595:G:H1	36:1:609:G:H5''	1.56	0.69
1:2:320:U:H2'	1:2:321:C:H2'	1.74	0.69
38:4:136:G:OP1	61:N5:48:SER:OG	2.06	0.69
1:6:1160:A:H2'	1:6:1161:C:C6	2.27	0.69
43:L6:175:LYS:O	50:M4:117:ARG:NH2	2.25	0.69
9:S7:78:THR:HA	9:S7:81:LEU:HB2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1097:U:H4'	1:6:1098:U:H5'	1.75	0.69
27:D5:74:SER:OG	1:6:1534:G:OP2	344.47	0.69
1:6:1600:A:H4'	1:6:1601:G:OP1	1.92	0.69
1:6:52:U:H2'	1:6:53:G:C8	2.28	0.69
6:S4:199:GLU:OE2	6:S4:209:HIS:NE2	2.24	0.69
8:S6:57:ASP:HA	8:S6:106:LEU:HA	1.74	0.69
42:L5:50:ARG:NH1	42:L5:72:ASP:OD2	2.25	0.69
44:L7:134:VAL:O	44:L7:229:PHE:HA	3.14	0.69
65:N9:23:LYS:HD2	65:N9:24:PRO:HD3	1.72	0.69
6:S4:199:GLU:HB2	6:S4:207:LEU:HB2	1.73	0.69
1:6:235:G:H2'	1:6:236:A:H8	1.57	0.69
15:C3:27:LYS:H	15:C3:27:LYS:HE3	1.57	0.69
20:C8:130:GLY:O	20:C8:145:ARG:NH1	2.25	0.69
55:M9:176:ARG:HA	55:M9:179:GLU:HB2	1.75	0.69
56:N0:6:GLU:OE1	56:N0:99:ARG:NH2	2.86	0.69
58:N2:56:VAL:HG13	58:N2:65:VAL:HG22	1.75	0.69
36:5:2537:U:O2'	36:5:2538:U:O4'	2.10	0.69
1:6:479:C:O2	1:6:510:G:N2	2.26	0.69
17:C5:30:THR:HG23	17:C5:86:VAL:HG21	1.73	0.69
47:M0:189:GLU:OE1	47:M0:202:LYS:NZ	2.21	0.69
64:N8:13:GLY:HA2	36:5:943:U:H3'	163.46	0.69
73:O7:14:LYS:HD3	75:O9:51:ILE:HG13	3.62	0.69
3:S1:150:VAL:HG23	1:6:1067:C:H5"	353.67	0.69
5:S3:117:ARG:HE	35:SM:122:GLU:HB3	1.57	0.69
18:C6:82:ARG:HH12	18:C6:114:ARG:HB3	1.56	0.69
46:L9:172:ILE:H	46:L9:172:ILE:HD13	1.58	0.69
46:L9:98:PRO:O	46:L9:116:ASN:ND2	2.25	0.69
73:O7:21:ARG:NH2	73:O7:41:ALA:O	2.30	0.69
4:S2:41:LEU:HD23	4:S2:240:LEU:HD11	2.23	0.69
1:2:1076:A:H4'	28:D6:13:LYS:HD3	1.74	0.69
36:5:1231:A:H5"	36:5:1232:C:H5'	1.75	0.69
13:C1:96:LYS:NZ	1:6:374:U:OP1	346.74	0.69
30:D8:42:ARG:HH22	30:D8:58:GLU:HG3	1.55	0.69
41:L4:287:THR:O	41:L4:291:ASN:ND2	2.32	0.69
42:L5:265:TYR:HE1	37:7:121:U:H5"	316.49	0.69
2:S0:30:GLN:HE22	2:S0:32:HIS:HB2	8.15	0.69
36:1:691:A:N1	38:4:28:C:O2'	2.20	0.69
1:2:1387:G:N7	19:C7:44:LYS:NZ	2.40	0.69
36:5:209:A:H4'	36:5:211:A:C8	2.27	0.69
1:6:228:G:N2	1:6:237:C:N3	2.41	0.69
1:2:1550:A:OP2	17:C5:42:ARG:NH2	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:M3:153:ASP:OD1	49:M3:157:ARG:NH2	2.25	0.69
6:S4:104:ASP:OD1	6:S4:104:ASP:N	2.26	0.69
34:SR:156:VAL:HA	34:SR:169:ILE:HG22	1.98	0.69
36:1:113:C:OP1	51:M5:147:ARG:NE	2.21	0.69
1:6:845:G:H2'	1:6:846:G:H8	1.56	0.69
33:E1:139:LEU:HD13	33:E1:151:ASN:HA	1.75	0.69
46:L9:126:VAL:HG21	46:L9:161:LEU:HA	1.75	0.69
47:M0:205:SER:OG	47:M0:208:ASN:OD1	3.44	0.69
36:1:595:G:N1	36:1:609:G:H5''	2.08	0.68
37:3:89:G:N2	37:3:92:A:OP2	2.26	0.68
21:C9:33:TYR:O	21:C9:35:ASP:N	3.74	0.68
25:D3:51:GLY:HA2	25:D3:77:ILE:HG13	1.86	0.68
4:S2:78:ASP:HB3	4:S2:104:VAL:HG12	4.80	0.68
36:5:2499:U:H2'	36:5:2500:A:H8	1.57	0.68
36:5:2812:C:H2'	36:5:2813:A:C8	2.27	0.68
1:6:1431:C:O2'	1:6:1437:U:O4	2.08	0.68
22:D0:69:LYS:HG2	22:D0:80:GLU:HB2	1.75	0.68
36:5:2714:G:N3	36:5:2714:G:H5''	2.08	0.68
36:1:2171:G:N2	39:L2:11:GLY:O	2.25	0.68
58:N2:54:VAL:HG12	58:N2:67:SER:HA	1.76	0.68
1:2:58:U:O2'	1:2:451:A:N3	2.25	0.68
1:2:535:A:OP2	11:S9:171:ARG:NH2	2.24	0.68
55:M9:5:ARG:NH1	36:5:1471:U:OP1	119.42	0.68
14:C2:67:THR:HG22	14:C2:68:GLU:HG3	1.74	0.68
1:2:458:G:OP1	26:D4:109:LYS:NZ	2.27	0.68
44:L7:105:LEU:O	36:5:1100:U:O2'	241.09	0.68
49:M3:129:ASN:HB2	49:M3:131:LYS:HE3	3.43	0.68
56:N0:13:ARG:HE	56:N0:51:VAL:HG22	4.80	0.68
71:O5:92:LEU:HB3	71:O5:96:GLU:HB2	1.75	0.68
20:C8:50:ALA:O	20:C8:68:ARG:NH1	4.86	0.68
63:N7:47:GLU:HB2	63:N7:69:LYS:HG2	4.50	0.68
16:C4:85:ALA:H	16:C4:119:THR:HG22	1.59	0.68
40:L3:37:ARG:HG2	40:L3:187:SER:H	4.26	0.68
44:L7:102:VAL:HG13	44:L7:126:LEU:HD23	1.74	0.68
48:M1:49:LYS:HB3	48:M1:62:ASN:HA	1.74	0.68
36:1:1389:G:H5''	68:O2:101:SER:HB3	1.75	0.68
79:Q3:46:THR:HB	79:Q3:58:SER:HB2	1.75	0.68
3:S1:171:ILE:HD12	3:S1:197:ILE:HD13	1.74	0.68
1:2:1561:U:H2'	1:2:1562:G:H8	1.58	0.68
1:2:1767:G:OP2	1:2:1770:U:O2'	2.11	0.68
36:5:2213:A:H2'	36:5:2214:A:C8	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:S8:2:GLY:N	1:6:393:C:OP2	291.49	0.68
1:6:75:U:O2'	1:6:76:A:O4'	2.12	0.68
40:L3:47:LEU:HG	40:L3:335:ILE:HD11	2.51	0.68
41:L4:158:SER:O	41:L4:160:GLN:NE2	2.27	0.68
47:M0:150:GLU:O	47:M0:153:ARG:N	4.21	0.68
49:M3:126:PHE:O	71:O5:114:ARG:NH2	2.57	0.68
55:M9:99:LEU:HD12	36:5:1722:U:H5''	224.39	0.68
71:O5:79:ASP:N	71:O5:79:ASP:OD1	2.88	0.68
3:S1:85:LYS:HB3	3:S1:101:HIS:HB3	2.88	0.68
36:1:3166:C:H42	36:1:3284:G:H1	1.41	0.68
43:L6:26:ARG:NH2	36:5:607:A:OP1	249.69	0.68
79:Q3:4:ARG:NH2	36:5:838:G:O6	236.28	0.68
1:6:1305:U:OP2	1:6:1306:C:N4	2.27	0.68
22:D0:23:ARG:NH2	1:6:1347:U:OP2	457.97	0.68
49:M3:56:PRO:HG3	49:M3:74:GLY:O	1.94	0.68
78:Q2:17:CYS:CB	78:Q2:77:CYS:SG	2.81	0.68
6:S4:117:GLU:O	6:S4:119:ALA:N	3.13	0.68
34:SR:180:ALA:HB3	34:SR:190:ALA:HB3	1.74	0.68
1:2:339:C:OP2	10:S8:10:LYS:NZ	2.26	0.68
38:4:84:C:O3'	62:N6:113:LYS:NZ	2.25	0.68
17:C5:15:HIS:H	17:C5:22:LEU:HD22	3.46	0.68
42:L5:83:LEU:HB3	42:L5:88:ILE:HB	1.95	0.68
50:M4:21:VAL:HG12	50:M4:65:LEU:HA	1.77	0.68
52:M6:36:VAL:HB	52:M6:108:ILE:HG22	4.72	0.68
66:O0:42:ILE:HG13	66:O0:67:VAL:HG13	1.76	0.68
3:S1:157:GLN:O	3:S1:159:SER:N	3.33	0.68
8:S6:10:ASN:ND2	8:S6:127:THR:O	3.52	0.68
1:6:884:A:H2'	1:6:885:G:C8	2.28	0.68
45:L8:43:LYS:HD2	61:N5:28:THR:HB	1.75	0.68
54:M8:38:ARG:NH2	36:5:1348:U:OP2	188.54	0.68
36:1:1786:G:H2'	36:1:1787:A:C8	2.29	0.67
64:N8:6:THR:HG22	64:N8:9:ARG:HG2	1.76	0.67
1:2:1251:U:H5'	33:E1:135:HIS:HD2	1.58	0.67
41:L4:60:THR:HG22	41:L4:62:ALA:H	3.01	0.67
47:M0:174:THR:OG1	47:M0:175:ASN:O	6.32	0.67
55:M9:11:ALA:HA	55:M9:41:ILE:HG21	2.33	0.67
63:N7:16:GLY:O	63:N7:18:TYR:N	2.28	0.67
70:O4:38:LEU:HD12	70:O4:38:LEU:H	2.83	0.67
9:S7:17:GLU:HG2	9:S7:46:ILE:HG22	1.76	0.67
11:S9:110:GLN:HE22	11:S9:126:ARG:HG2	1.58	0.67
11:S9:27:GLU:HB3	11:S9:39:LYS:HD2	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:284:A:OP2	78:Q2:41:ARG:NH1	2.27	0.67
1:2:885:G:N2	1:2:927:C:O2	2.23	0.67
8:S6:159:ARG:NH2	1:6:79:C:OP1	349.40	0.67
16:C4:25:ASP:OD1	16:C4:26:THR:N	3.06	0.67
25:D3:70:LYS:HB3	25:D3:93:LEU:HD22	1.76	0.67
3:S1:48:VAL:HG21	3:S1:61:LEU:HD22	5.24	0.67
4:S2:37:PRO:HG3	4:S2:46:LYS:HD2	3.49	0.67
11:S9:61:THR:HA	24:D2:97:ARG:HH12	1.64	0.67
36:1:751:A:H2'	36:1:752:C:H6	1.60	0.67
56:N0:137:ARG:NH2	36:5:1214:U:OP2	329.98	0.67
15:C3:127:ARG:NH2	1:6:629:U:OP1	308.01	0.67
1:6:891:A:H2'	1:6:892:A:C8	2.29	0.67
40:L3:50:LYS:HG2	40:L3:332:ARG:HA	1.76	0.67
41:L4:3:ARG:O	41:L4:5:GLN:N	2.28	0.67
50:M4:16:GLU:HB3	56:N0:149:LYS:HB3	1.99	0.67
57:N1:78:LYS:HB3	57:N1:87:LYS:HG3	1.77	0.67
77:Q1:2:ARG:NH1	1:6:1773:C:OP2	309.23	0.67
6:S4:180:LEU:HA	6:S4:194:THR:HA	1.77	0.67
36:1:2759:U:H5''	36:1:2760:C:H5'	1.75	0.67
36:1:542:G:H2'	36:1:543:C:C6	2.29	0.67
1:2:1253:U:H5''	33:E1:130:VAL:HB	1.76	0.67
1:2:145:A:O2'	1:2:146:U:O5'	2.12	0.67
20:C8:13:HIS:HA	20:C8:24:GLY:HA3	1.75	0.67
26:D4:124:ARG:HA	26:D4:127:LYS:HG2	1.76	0.67
49:M3:108:ILE:O	49:M3:112:ASN:ND2	2.27	0.67
50:M4:121:MET:HE1	36:5:3215:A:H8	277.43	0.67
6:S4:18:TRP:O	6:S4:51:ARG:NH2	4.46	0.67
35:SM:43:ASP:OD1	35:SM:45:SER:OG	3.08	0.67
25:D3:109:ARG:HB3	25:D3:112:LYS:HB2	1.77	0.67
39:L2:193:ARG:NH1	36:5:2174:G:OP2	190.72	0.67
41:L4:141:ARG:O	41:L4:143:GLU:N	4.55	0.67
41:L4:338:LYS:O	41:L4:340:GLY:N	2.28	0.67
48:M1:137:ARG:NH2	37:7:44:C:OP2	296.85	0.67
68:O2:69:SER:OG	68:O2:70:GLY:N	2.24	0.67
36:5:997:A:O2'	37:7:79:A:N3	2.28	0.67
2:S0:184:LEU:HD12	23:D1:45:ALA:HB2	1.77	0.67
41:L4:141:ARG:C	41:L4:143:GLU:H	3.74	0.67
41:L4:354:VAL:HG11	57:N1:143:THR:HG21	2.83	0.67
49:M3:46:ILE:HG22	49:M3:49:ARG:HB3	1.77	0.67
53:M7:111:LYS:HA	53:M7:153:LYS:HE2	4.52	0.67
58:N2:82:LYS:NZ	36:5:1682:U:O2	160.01	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
71:O5:85:THR:HG22	71:O5:88:LEU:H	2.07	0.67
36:1:3067:C:OP1	55:M9:58:HIS:NE2	2.19	0.67
26:D4:83:LYS:HE2	26:D4:96:LEU:HB3	1.76	0.67
44:L7:103:LEU:HA	44:L7:130:ILE:HD11	4.72	0.67
36:1:900:G:H1'	36:1:1589:A:N6	2.10	0.67
36:1:3094:A:H2'	36:1:3095:U:C6	2.30	0.67
36:5:900:G:H1'	36:5:1589:A:N6	2.10	0.67
33:E1:146:SER:HB3	1:6:1234:A:H4'	434.46	0.67
45:L8:115:ALA:O	45:L8:117:ALA:N	2.26	0.67
36:1:976:U:OP1	54:M8:144:ARG:NH2	2.27	0.67
11:S9:11:THR:HG23	1:6:472:U:H5''	397.66	0.67
11:S9:109:LEU:HB2	11:S9:146:PHE:HB3	1.79	0.67
36:1:655:C:H2'	36:1:656:A:H8	1.58	0.67
1:2:1235:C:H5'	33:E1:146:SER:HB2	1.77	0.67
44:L7:98:LYS:HD2	36:5:985:U:H5''	242.12	0.67
18:C6:13:LYS:HD2	18:C6:76:SER:HA	1.77	0.67
24:D2:94:LEU:HD11	24:D2:102:VAL:HG23	2.07	0.67
40:L3:147:GLU:OE2	40:L3:150:ARG:NH1	2.89	0.67
51:M5:147:ARG:NH2	36:5:113:C:OP1	77.60	0.67
52:M6:65:ASN:ND2	36:5:2988:C:OP1	220.91	0.67
63:N7:88:ASP:HB3	63:N7:121:ARG:HH22	1.60	0.67
36:1:801:A:OP1	64:N8:27:LYS:NZ	2.27	0.67
64:N8:96:LYS:O	64:N8:98:THR:N	2.27	0.67
9:S7:73:VAL:O	9:S7:75:THR:N	2.28	0.67
1:2:853:G:N7	55:M9:173:ARG:NH2	2.43	0.66
38:4:22:U:OP1	62:N6:12:ARG:NH2	2.26	0.66
1:6:67:A:O2'	1:6:69:G:OP1	2.09	0.66
1:6:800:U:H2'	1:6:801:G:C8	2.30	0.66
25:D3:68:ILE:O	25:D3:70:LYS:NZ	2.28	0.66
51:M5:49:ARG:NH1	36:5:149:U:OP2	100.65	0.66
36:1:994:G:H3'	57:N1:13:TYR:CD2	2.30	0.66
61:N5:47:ALA:HB3	71:O5:77:PRO:HG3	1.78	0.66
76:Q0:77:ILE:HG23	76:Q0:78:ILE:HG22	9.14	0.66
3:S1:129:THR:OG1	3:S1:131:ASP:O	2.87	0.66
3:S1:131:ASP:O	3:S1:133:TYR:N	2.28	0.66
36:1:2660:G:O3'	36:1:2749:G:N2	2.28	0.66
1:2:1018:U:H2'	1:2:1019:A:H8	1.60	0.66
1:2:2:A:O2'	4:S2:198:THR:O	2.12	0.66
1:2:778:G:H2'	1:2:779:U:H2'	1.76	0.66
13:C1:124:THR:HB	13:C1:141:LYS:HB3	2.10	0.66
41:L4:35:VAL:HG21	41:L4:244:LEU:HD21	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S2:164:SER:OG	4:S2:165:VAL:N	2.95	0.66
36:1:2193:U:H5'	36:1:2194:G:H5'	1.75	0.66
36:1:2687:G:OP1	42:L5:8:LYS:NZ	2.27	0.66
51:M5:49:ARG:NH2	36:5:115:A:OP1	100.96	0.66
32:E0:37:ARG:NH1	1:6:478:A:OP1	439.99	0.66
1:6:778:G:OP2	1:6:780:A:N6	2.28	0.66
22:D0:21:LYS:HB3	22:D0:94:GLU:HG2	4.47	0.66
52:M6:3:VAL:HG13	52:M6:4:GLU:HG3	1.77	0.66
6:S4:124:GLY:HA2	6:S4:142:HIS:CE1	2.30	0.66
36:5:618:C:O2'	36:5:621:A:N3	2.23	0.66
1:6:800:U:H2'	1:6:801:G:H8	1.60	0.66
20:C8:120:ARG:NH2	35:SM:58:GLU:OE1	2.28	0.66
44:L7:110:ARG:HD2	54:M8:3:ILE:HG23	4.01	0.66
61:N5:100:LYS:HE3	61:N5:106:ASP:HA	1.76	0.66
36:1:1598:G:OP2	70:O4:31:ARG:NH2	2.29	0.66
36:1:1613:A:OP2	74:O8:46:ARG:NH2	2.28	0.66
10:S8:61:GLU:HG2	10:S8:62:THR:HG23	2.97	0.66
1:6:778:G:N2	1:6:780:A:H5'	2.11	0.66
45:L8:162:LEU:HA	51:M5:7:LEU:HD11	1.77	0.66
2:S0:140:ASN:ND2	4:S2:60:SER:O	2.29	0.66
36:1:352:A:H61	36:1:365:A:H5''	1.60	0.66
47:M0:63:GLU:HB2	36:5:2853:A:H5'	296.05	0.66
64:N8:36:GLY:HA3	64:N8:40:HIS:CE1	2.30	0.66
2:S0:134:LYS:O	2:S0:137:SER:OG	2.11	0.66
11:S9:3:ARG:H	11:S9:3:ARG:HH21	1.41	0.66
36:1:1665:C:H2'	36:1:1666:G:H8	1.59	0.66
36:1:528:U:H2'	36:1:529:A:C8	2.30	0.66
1:6:1039:A:O2'	1:6:1040:G:O5'	2.14	0.66
1:6:907:A:N3	1:6:997:G:O2'	2.23	0.66
15:C3:136:PRO:HG2	15:C3:139:TRP:HB2	1.77	0.66
23:D1:3:ASN:HD21	23:D1:7:GLN:HB3	1.61	0.66
4:S2:47:ALA:O	4:S2:49:LYS:N	2.29	0.66
1:2:1657:U:H4'	1:2:1658:G:O5'	1.96	0.66
1:6:1711:C:H2'	1:6:1712:A:H5''	1.78	0.66
10:S8:16:ALA:HB2	1:6:354:C:H5''	297.62	0.66
23:D1:21:ASN:OD1	24:D2:23:ARG:NH2	2.27	0.66
40:L3:3:HIS:O	40:L3:5:LYS:N	2.28	0.66
46:L9:171:ASP:OD2	46:L9:173:ARG:NH1	2.28	0.66
59:N3:101:VAL:HG11	59:N3:114:ILE:HG12	1.78	0.66
59:N3:6:ALA:HB1	59:N3:125:LEU:HD11	1.76	0.66
62:N6:27:ARG:HA	62:N6:30:LEU:HD12	2.49	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:14:VAL:HG12	63:N7:15:ARG:HG3	1.76	0.66
4:S2:163:GLY:O	4:S2:165:VAL:N	4.24	0.66
5:S3:141:LYS:HE2	5:S3:180:GLY:HA3	1.78	0.66
9:S7:83:LYS:HE3	9:S7:84:LYS:HE2	1.77	0.66
36:1:1230:G:H1	36:1:1279:C:H42	1.44	0.66
36:1:1355:A:H5''	36:1:1356:U:H5	1.61	0.66
36:1:2228:A:H2'	36:1:2229:A:C8	2.31	0.66
36:5:1103:A:H5''	36:5:1104:G:OP2	1.95	0.66
1:6:485:A:N6	1:6:486:G:N3	2.44	0.66
1:6:895:G:H1	1:6:917:U:H3	1.43	0.66
22:D0:41:ILE:HG13	22:D0:107:THR:HG21	4.17	0.66
39:L2:206:PRO:HG3	39:L2:213:GLY:HA2	3.02	0.66
40:L3:56:ILE:HD13	40:L3:76:VAL:HG21	1.76	0.66
47:M0:170:LYS:HA	47:M0:177:ASP:HA	2.09	0.66
56:N0:135:VAL:O	56:N0:141:LYS:NZ	2.22	0.66
34:SR:13:LEU:HB2	34:SR:310:ILE:HB	1.77	0.66
36:1:2261:G:O2'	36:1:2263:C:N4	2.29	0.66
1:2:115:G:OP1	13:C1:67:ARG:NH1	2.29	0.66
16:C4:26:THR:HG21	16:C4:97:GLY:HA3	2.07	0.66
19:C7:28:PHE:HA	19:C7:55:THR:HG21	3.64	0.66
26:D4:61:ARG:NH2	1:6:530:C:O2	409.02	0.66
30:D8:15:VAL:HA	30:D8:28:VAL:HG22	1.76	0.66
49:M3:166:ALA:N	64:N8:135:GLU:OE1	2.27	0.66
53:M7:27:LYS:HD3	53:M7:63:PHE:HB3	1.78	0.66
63:N7:26:VAL:HG21	63:N7:96:VAL:HB	1.78	0.66
36:1:2895:G:O2'	76:Q0:100:TYR:O	2.13	0.66
1:2:7:G:N7	4:S2:205:ARG:NH1	2.43	0.66
7:S5:124:LEU:O	27:D5:58:ARG:NH1	6.39	0.66
10:S8:192:TYR:O	10:S8:196:LEU:HB2	1.96	0.66
11:S9:54:ARG:HB2	11:S9:57:ARG:HH21	3.22	0.66
36:5:2102:U:H2'	36:5:2103:U:H6	1.61	0.65
12:C0:7:ASP:OD2	12:C0:10:LYS:NZ	4.95	0.65
36:1:911:C:H5''	39:L2:15:ILE:HD13	1.78	0.65
36:1:2688:U:OP1	42:L5:12:TYR:OH	2.12	0.65
55:M9:175:GLN:HG2	55:M9:176:ARG:HH21	1.62	0.65
2:S0:108:THR:OG1	2:S0:135:GLU:OE1	2.65	0.65
1:2:1266:U:H2'	1:2:1267:G:H8	1.61	0.65
1:2:169:A:H5''	8:S6:176:GLN:HG2	1.79	0.65
15:C3:115:LEU:HD22	15:C3:119:GLU:HG3	2.26	0.65
25:D3:56:LYS:NZ	25:D3:96:VAL:O	5.67	0.65
1:2:1234:A:H4'	33:E1:146:SER:HB3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:M7:116:HIS:HB3	53:M7:149:VAL:HB	2.07	0.65
5:S3:70:THR:HG22	5:S3:86:LEU:HB2	2.14	0.65
10:S8:36:THR:HG21	10:S8:173:PRO:HB2	1.78	0.65
36:5:1355:A:H5''	36:5:1357:G:H1'	1.79	0.65
2:S0:200:ASP:HB2	19:C7:85:VAL:HG22	1.77	0.65
22:D0:31:VAL:HA	22:D0:34:LEU:HB3	2.80	0.65
25:D3:90:ASP:O	25:D3:136:TRP:NE1	2.29	0.65
26:D4:52:LYS:O	26:D4:54:ALA:N	2.37	0.65
43:L6:40:LEU:HD13	43:L6:84:VAL:HG11	2.24	0.65
46:L9:101:VAL:HG22	46:L9:114:VAL:HG22	1.79	0.65
58:N2:19:VAL:HG12	58:N2:105:LEU:HD22	1.99	0.65
62:N6:99:LEU:HD13	62:N6:104:LEU:HD21	1.78	0.65
68:O2:19:ARG:HE	68:O2:33:ARG:HB2	2.07	0.65
1:2:1728:A:H1'	10:S8:32:GLN:NE2	2.12	0.65
1:2:1797:A:OP2	28:D6:10:ARG:NH2	2.30	0.65
1:2:72:A:N7	8:S6:169:TYR:OH	2.25	0.65
36:5:1717:U:H2'	36:5:1718:G:C8	2.32	0.65
36:5:1940:G:H21	36:5:3362:A:H8	1.44	0.65
17:C5:86:VAL:O	17:C5:88:GLU:N	2.29	0.65
32:E0:12:GLY:O	32:E0:16:SER:OG	2.89	0.65
42:L5:40:HIS:CE1	42:L5:42:ALA:HB3	3.72	0.65
56:N0:13:ARG:HB3	56:N0:51:VAL:HG11	1.79	0.65
7:S5:90:ILE:HD11	7:S5:130:ILE:HG13	1.84	0.65
9:S7:151:LYS:HA	9:S7:182:VAL:HG12	1.77	0.65
1:6:1087:A:H2'	1:6:1088:A:C8	2.32	0.65
45:L8:45:ASN:HD21	45:L8:47:SER:HB3	1.62	0.65
43:L6:51:ARG:NH1	50:M4:114:ASP:OD2	2.30	0.65
57:N1:14:MET:HE1	57:N1:55:LYS:HB2	2.34	0.65
3:S1:157:GLN:HB2	3:S1:160:HIS:CE1	3.42	0.65
20:C8:144:ARG:O	35:SM:68:ARG:NH2	2.30	0.65
34:SR:42:LEU:HD21	34:SR:82:SER:HB3	2.04	0.65
28:D6:15:ARG:NH1	1:6:936:G:N7	319.07	0.65
15:C3:66:ILE:HG13	15:C3:67:THR:HG22	2.78	0.65
37:3:46:A:OP2	42:L5:158:ARG:NH1	2.30	0.65
49:M3:91:ARG:NH2	49:M3:97:VAL:O	2.24	0.65
36:1:2093:A:N1	55:M9:114:LYS:NZ	2.45	0.65
55:M9:23:TRP:HB3	55:M9:51:VAL:HG22	1.79	0.65
61:N5:34:LEU:HD22	61:N5:35:PRO:HD2	1.99	0.65
75:O9:28:ARG:HH11	75:O9:36:ARG:HD3	6.82	0.65
75:O9:43:ASN:HB3	75:O9:46:ARG:HG3	4.01	0.65
11:S9:93:LEU:HA	11:S9:96:VAL:HG13	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:2557:A:OP1	39:L2:69:TYR:OH	2.14	0.65
1:2:551:G:H1	1:2:573:C:H42	1.45	0.65
36:5:3194:C:H2'	36:5:3195:U:H3'	1.79	0.65
27:D5:66:VAL:HA	27:D5:71:ILE:HG22	7.24	0.65
41:L4:309:ARG:HH21	41:L4:312:VAL:HB	2.97	0.65
49:M3:28:GLN:OE1	51:M5:201:ARG:NH1	2.28	0.65
51:M5:47:LYS:HA	51:M5:50:ARG:HE	1.61	0.65
1:2:1126:G:OP2	77:Q1:18:ARG:NH2	2.29	0.65
2:S0:157:ASP:OD2	23:D1:60:ARG:NH2	2.30	0.65
8:S6:2:LYS:HB3	8:S6:108:VAL:HG12	5.64	0.65
36:1:239:G:O2'	36:1:240:U:OP1	2.15	0.65
36:1:685:G:OP2	49:M3:35:ARG:NH1	2.29	0.65
1:2:1344:A:N6	1:2:1377:U:O2'	2.29	0.65
1:2:190:C:N4	1:2:196:G:O6	2.30	0.65
38:8:26:U:H2'	38:8:27:U:C6	2.32	0.65
1:2:1258:U:H4'	12:C0:2:LEU:HD13	1.79	0.65
21:C9:39:THR:HA	21:C9:100:ILE:HD12	4.12	0.65
40:L3:84:VAL:HG22	40:L3:162:VAL:HB	2.90	0.65
49:M3:47:ALA:HB1	49:M3:48:PRO:HD2	1.79	0.65
51:M5:115:VAL:HG22	51:M5:134:LEU:HD23	1.79	0.65
55:M9:106:LEU:HD13	55:M9:138:LEU:HD11	2.56	0.65
36:1:2738:A:H5''	65:N9:38:LYS:HG3	1.79	0.65
36:5:1506:A:H1'	36:5:1848:G:O6	1.97	0.65
1:6:138:A:H62	1:6:266:A:H61	1.43	0.65
16:C4:107:ARG:NH2	28:D6:52:ASP:OD2	4.11	0.65
24:D2:18:GLU:HG3	24:D2:69:LEU:HD23	1.79	0.65
39:L2:144:ASN:ND2	39:L2:161:ASP:OD2	3.74	0.65
40:L3:205:VAL:HG11	40:L3:322:ILE:HD11	2.44	0.65
45:L8:156:ASP:OD2	45:L8:156:ASP:N	2.28	0.65
46:L9:105:GLU:HG3	46:L9:109:ALA:HB3	1.78	0.65
48:M1:90:GLN:HG2	48:M1:170:ASP:HB2	1.79	0.65
53:M7:109:ALA:HA	53:M7:112:LEU:HD22	1.94	0.65
63:N7:29:HIS:O	63:N7:31:GLU:N	2.30	0.65
73:O7:31:LYS:NZ	36:5:815:G:OP2	136.34	0.65
4:S2:227:PRO:HA	4:S2:230:TRP:CE2	2.32	0.65
5:S3:80:ALA:O	5:S3:83:THR:OG1	2.27	0.65
7:S5:94:THR:HG22	7:S5:114:ILE:HG13	1.89	0.65
36:5:2696:A:H2'	36:5:2697:A:C8	2.31	0.65
36:5:980:A:H2'	36:5:981:U:C2	2.32	0.65
1:6:489:C:O2'	1:6:490:C:O4'	2.14	0.65
1:6:513:U:H2'	1:6:514:G:C8	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:C0:4:PRO:HG2	12:C0:7:ASP:HB2	2.96	0.65
19:C7:66:VAL:O	19:C7:68:GLY:N	3.66	0.65
41:L4:193:LYS:HB2	41:L4:198:ARG:HA	3.33	0.65
47:M0:14:ASN:O	47:M0:128:ARG:NH2	2.43	0.65
51:M5:110:ALA:HB1	51:M5:113:LEU:HD23	1.79	0.65
67:O1:75:ILE:HG12	67:O1:93:VAL:HG13	2.67	0.65
5:S3:40:ARG:HG2	22:D0:110:PRO:HB3	1.93	0.65
6:S4:66:MET:SD	6:S4:78:THR:OG1	2.86	0.65
36:1:2307:G:O2'	36:1:2310:U:OP2	2.15	0.64
1:2:751:G:H2'	1:2:752:A:H8	1.62	0.64
28:D6:76:SER:OG	1:6:1793:G:N2	320.92	0.64
41:L4:55:LYS:HE2	41:L4:59:GLN:HE22	1.62	0.64
53:M7:52:LEU:HD21	53:M7:89:LYS:HD2	5.31	0.64
57:N1:12:ARG:HD2	57:N1:13:TYR:CE1	2.32	0.64
71:O5:38:ARG:HH11	71:O5:41:LEU:HD13	1.61	0.64
75:O9:17:LYS:NZ	36:5:1521:G:O6	96.75	0.64
6:S4:125:LYS:HZ3	6:S4:157:ASN:HA	5.15	0.64
6:S4:59:ARG:NH2	1:6:445:A:OP2	382.39	0.64
7:S5:157:ARG:NH1	7:S5:224:ASN:OD1	3.82	0.64
34:SR:89:LEU:HB2	34:SR:103:PHE:HB2	1.78	0.64
34:SR:80:ALA:HB3	34:SR:92:TRP:HB2	1.95	0.64
1:2:1242:A:OP1	17:C5:59:LYS:NZ	2.28	0.64
1:2:647:G:H2'	1:2:648:G:H8	1.62	0.64
1:6:319:U:H5''	1:6:320:U:H5	1.62	0.64
14:C2:63:VAL:HG11	14:C2:94:ALA:HA	1.80	0.64
29:D7:67:THR:OG1	29:D7:70:LYS:O	2.15	0.64
36:1:2146:C:OP1	39:L2:200:ARG:NH1	2.30	0.64
40:L3:76:VAL:HG12	40:L3:325:LYS:HA	1.78	0.64
71:O5:81:ARG:NH2	36:5:18:G:OP1	76.41	0.64
36:5:1024:G:O6	36:5:1029:G:N2	2.29	0.64
45:L8:240:ASN:ND2	36:5:2586:G:OP1	181.76	0.64
47:M0:158:LYS:NZ	36:5:2852:C:N3	307.76	0.64
1:6:1594:G:OP2	1:6:1596:C:N4	2.30	0.64
1:6:1603:U:H2'	1:6:1604:U:H6	1.62	0.64
1:6:51:A:OP2	1:6:424:C:N4	2.27	0.64
1:6:653:C:H42	1:6:677:G:H1	1.46	0.64
23:D1:64:GLU:HG3	29:D7:3:LEU:HG	3.09	0.64
30:D8:13:ILE:HB	30:D8:29:ARG:HG2	5.50	0.64
42:L5:40:HIS:HB3	42:L5:43:LYS:HG3	1.79	0.64
50:M4:135:LEU:HD22	52:M6:177:LYS:HE2	1.79	0.64
51:M5:94:TYR:CE1	51:M5:96:ARG:HB2	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
61:N5:100:LYS:NZ	61:N5:106:ASP:OD2	5.49	0.64
74:O8:9:LYS:NZ	74:O8:13:GLU:OE1	2.25	0.64
4:S2:45:VAL:HG21	4:S2:68:ILE:HG23	1.79	0.64
36:1:873:C:H5''	36:1:874:U:O5'	1.98	0.64
1:2:1696:G:H21	1:2:1705:C:H41	1.45	0.64
1:2:738:G:H2'	1:2:739:G:H8	1.63	0.64
36:5:503:C:H2'	36:5:504:A:C8	2.32	0.64
1:6:538:A:N7	1:6:543:C:N4	2.40	0.64
20:C8:52:VAL:HG13	20:C8:61:LEU:HD21	3.11	0.64
24:D2:77:PRO:O	24:D2:79:PHE:N	2.44	0.64
26:D4:112:LYS:NZ	26:D4:113:ASN:OD1	2.90	0.64
45:L8:221:ASN:HA	45:L8:225:LYS:HE3	2.15	0.64
11:S9:74:ASN:HA	11:S9:77:ILE:HD12	1.78	0.64
1:2:654:C:H3'	1:2:655:G:H5''	1.78	0.64
57:N1:49:GLN:HG2	36:5:2756:C:O4'	247.07	0.64
36:5:437:G:H1	36:5:622:A:H61	1.44	0.64
33:E1:144:CYS:HB3	33:E1:147:VAL:HG22	1.79	0.64
62:N6:48:LEU:HD23	62:N6:49:PRO:HD2	2.46	0.64
11:S9:103:ASP:OD2	11:S9:103:ASP:N	2.83	0.64
1:2:1280:C:H2'	1:2:1281:G:H8	1.62	0.64
1:2:1794:A:H1'	28:D6:79:ILE:HD12	1.79	0.64
1:6:27:U:H2'	1:6:28:A:H8	1.63	0.64
16:C4:97:GLY:HA2	16:C4:100:ALA:HB3	1.79	0.64
5:S3:11:LEU:HD12	22:D0:86:ILE:HG12	1.80	0.64
36:1:598:A:OP1	44:L7:41:ARG:NH1	2.30	0.64
46:L9:27:VAL:HG11	46:L9:79:ILE:HA	1.80	0.64
48:M1:37:LEU:O	48:M1:41:SER:OG	2.09	0.64
53:M7:18:ARG:NH2	53:M7:147:GLU:OE1	2.30	0.64
63:N7:36:HIS:HD2	63:N7:74:VAL:HG11	3.84	0.64
69:O3:42:GLN:HA	69:O3:45:LEU:HD23	4.28	0.64
6:S4:126:VAL:HA	6:S4:141:THR:HA	1.80	0.64
36:5:655:C:H2'	36:5:656:A:C8	2.33	0.64
1:6:591:A:H2'	1:6:592:A:H8	1.61	0.64
17:C5:129:GLY:HA3	35:SM:74:LYS:HD2	5.58	0.64
22:D0:50:LEU:HD21	22:D0:95:ALA:HB2	1.78	0.64
39:L2:70:ARG:HH21	39:L2:72:ARG:HH21	6.36	0.64
40:L3:260:VAL:HG11	40:L3:266:ARG:HH11	1.62	0.64
55:M9:105:LEU:HD12	55:M9:135:LYS:HG3	1.78	0.64
34:SR:12:THR:OG1	34:SR:14:GLU:OE1	6.50	0.64
29:D7:70:LYS:HG3	1:6:1049:U:H5''	347.60	0.64
25:D3:65:ASN:ND2	1:6:574:G:O6	364.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:C1:6:THR:O	13:C1:8:GLN:N	2.31	0.64
16:C4:43:THR:OG1	16:C4:44:GLY:N	2.31	0.64
23:D1:81:ASN:O	23:D1:83:TRP:N	2.31	0.64
43:L6:40:LEU:HD12	43:L6:65:ILE:HD11	6.56	0.64
49:M3:54:LEU:N	49:M3:94:GLY:O	2.67	0.64
54:M8:122:ILE:HG23	54:M8:126:GLN:HB2	1.80	0.64
56:N0:1:MET:HE1	56:N0:32:SER:H	1.62	0.64
72:O6:4:LYS:HD2	72:O6:14:GLY:HA3	2.76	0.64
78:Q2:17:CYS:SG	78:Q2:77:CYS:CB	2.86	0.64
36:5:1750:A:H4'	36:5:1751:G:H5'	1.79	0.64
71:O5:90:ARG:NH1	36:5:20:A:OP2	86.01	0.64
13:C1:109:VAL:HG12	13:C1:137:PHE:HB2	4.84	0.64
28:D6:88:SER:OG	28:D6:90:GLU:OE2	6.12	0.64
40:L3:10:ARG:NH2	40:L3:263:SER:O	2.31	0.64
50:M4:120:VAL:HG22	52:M6:197:LEU:HD13	1.80	0.64
51:M5:42:PRO:HG3	51:M5:61:ILE:HG13	2.40	0.64
52:M6:22:VAL:HG11	52:M6:120:VAL:HG11	1.80	0.64
58:N2:67:SER:OG	58:N2:69:ALA:O	3.75	0.64
60:N4:36:SER:O	60:N4:40:PHE:N	2.31	0.64
36:5:251:G:N7	36:5:253:A:N6	2.46	0.64
14:C2:59:LEU:HA	14:C2:87:PRO:HB2	1.80	0.64
1:2:576:G:O6	25:D3:65:ASN:ND2	2.31	0.64
27:D5:41:ILE:HG13	27:D5:42:LEU:H	1.62	0.64
42:L5:216:GLU:O	42:L5:220:SER:OG	2.11	0.64
43:L6:40:LEU:HB2	43:L6:52:VAL:HG12	2.07	0.64
50:M4:45:LEU:HD22	56:N0:72:VAL:HG23	2.64	0.64
34:SR:295:SER:HB3	34:SR:302:PHE:HE2	3.47	0.64
36:1:541:U:H2'	36:1:542:G:C8	2.33	0.63
1:2:1500:C:OP1	21:C9:122:ARG:NH2	2.27	0.63
46:L9:116:ASN:O	46:L9:120:ASP:N	2.30	0.63
61:N5:105:VAL:HG13	61:N5:130:TYR:CD2	2.46	0.63
64:N8:74:ASN:HB2	64:N8:76:ASP:HB2	1.80	0.63
67:O1:72:ARG:NH1	67:O1:105:GLN:O	2.30	0.63
73:O7:19:CYS:SG	73:O7:34:CYS:HB2	2.37	0.63
4:S2:137:ILE:HG12	4:S2:138:PRO:HD2	1.80	0.63
36:1:1068:C:OP1	57:N1:110:LYS:NZ	2.31	0.63
36:1:1550:C:H2'	36:1:1551:C:H6	1.64	0.63
36:5:3295:A:H2'	36:5:3296:A:C8	2.34	0.63
20:C8:87:ASN:OD1	20:C8:88:ARG:N	2.30	0.63
3:S1:225:VAL:HA	3:S1:228:LEU:HB3	2.94	0.63
11:S9:163:PRO:HB3	11:S9:169:PRO:HA	3.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1101:G:OP2	44:L7:196:LYS:NZ	2.27	0.63
36:1:1242:G:N2	36:1:1270:A:O2'	2.31	0.63
36:1:979:U:H1'	36:1:980:A:C5	2.34	0.63
1:2:1280:C:H2'	1:2:1281:G:C8	2.33	0.63
1:2:1534:G:N7	27:D5:77:ARG:NH2	2.46	0.63
36:5:348:A:N3	36:5:352:A:O2'	2.31	0.63
21:C9:57:ARG:NH1	1:6:1479:A:OP1	391.42	0.63
6:S4:69:HIS:HB3	26:D4:17:LEU:HD23	1.80	0.63
42:L5:160:PHE:CD2	42:L5:179:ARG:HB3	2.54	0.63
51:M5:59:PHE:HD1	51:M5:133:ILE:HD11	1.63	0.63
52:M6:77:SER:OG	52:M6:106:GLU:OE1	2.13	0.63
5:S3:141:LYS:NZ	5:S3:179:GLN:OE1	4.79	0.63
1:2:229:U:H3	1:2:236:A:H61	1.47	0.63
1:6:1163:A:N3	1:6:1613:U:O2'	2.26	0.63
1:2:1553:G:H4'	31:D9:14:TYR:HE1	1.62	0.63
46:L9:91:ARG:HG2	46:L9:182:SER:HB3	5.98	0.63
62:N6:37:LYS:H	62:N6:37:LYS:HD3	1.63	0.63
36:1:965:A:H2	64:N8:43:ILE:HD12	1.61	0.63
2:S0:29:VAL:HG12	2:S0:149:LEU:HB3	7.42	0.63
8:S6:163:THR:HB	8:S6:168:THR:HG22	3.74	0.63
36:1:915:A:H8	36:1:2136:C:HO2'	1.47	0.63
51:M5:69:GLY:O	36:5:290:G:H4'	145.47	0.63
21:C9:125:SER:O	21:C9:129:GLN:N	2.60	0.63
23:D1:71:ARG:HG3	29:D7:4:VAL:HG11	1.80	0.63
46:L9:163:GLN:O	46:L9:166:ARG:HD3	1.98	0.63
46:L9:77:ASN:HA	46:L9:80:THR:HG23	3.95	0.63
55:M9:127:SER:HB3	55:M9:132:PHE:HD2	3.37	0.63
36:1:1864:A:OP1	55:M9:88:ARG:NH1	2.31	0.63
72:O6:54:GLU:HG2	72:O6:90:MET:HE1	2.53	0.63
36:1:516:A:O3'	44:L7:60:ARG:NH2	2.25	0.63
13:C1:22:ASN:HD21	13:C1:24:LYS:HB2	1.63	0.63
26:D4:33:ALA:O	26:D4:34:ASN:ND2	2.29	0.63
36:1:2852:C:N3	47:M0:158:LYS:NZ	2.46	0.63
48:M1:37:LEU:HD12	48:M1:67:VAL:HG13	4.99	0.63
61:N5:80:ASN:ND2	61:N5:126:LEU:O	2.32	0.63
63:N7:25:ILE:HA	63:N7:43:VAL:HG12	1.79	0.63
9:S7:75:THR:HG22	9:S7:161:GLN:HE22	4.20	0.63
10:S8:39:GLY:HA2	10:S8:61:GLU:HB3	1.80	0.63
10:S8:98:LYS:HB3	1:6:329:G:H5''	274.27	0.63
36:1:415:G:H2'	36:1:416:A:C8	2.34	0.63
36:1:715:A:HO2'	36:1:752:C:HO2'	1.46	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:L2:243:THR:HG23	36:5:2242:A:H5'	232.22	0.63
18:C6:34:SER:HB3	18:C6:38:LEU:HD12	1.81	0.63
9:S7:144:VAL:HG22	24:D2:49:GLU:HB2	5.40	0.63
43:L6:40:LEU:N	43:L6:52:VAL:O	2.65	0.63
52:M6:61:ALA:HA	52:M6:70:PRO:HD2	1.81	0.63
71:O5:21:LEU:HD21	71:O5:51:ILE:HG23	3.11	0.63
2:S0:126:PRO:HB2	2:S0:152:PRO:HG2	2.07	0.63
2:S0:133:ILE:HG23	2:S0:155:PHE:HB2	1.81	0.63
36:1:2623:G:H2'	36:1:2624:G:H8	1.63	0.63
36:1:903:U:OP2	73:O7:30:GLN:NE2	2.31	0.63
1:2:479:C:O2	1:2:510:G:N2	2.28	0.63
36:5:1103:A:H3'	36:5:1104:G:H5'	1.79	0.63
45:L8:38:GLN:HB2	36:5:2557:A:H2	207.06	0.63
36:5:2592:G:H4'	36:5:2594:C:C2	2.34	0.63
36:5:3280:U:O2'	36:5:3281:U:H5''	1.98	0.63
16:C4:29:HIS:HB3	16:C4:41:ARG:HG3	1.80	0.63
41:L4:6:VAL:N	41:L4:20:LEU:O	2.43	0.63
47:M0:36:LEU:HD21	47:M0:69:ARG:HD3	2.07	0.63
62:N6:58:VAL:HA	62:N6:104:LEU:HD23	1.79	0.63
70:O4:67:LYS:HA	70:O4:70:LYS:HE3	1.81	0.63
2:S0:160:ILE:O	2:S0:162:CYS:N	2.31	0.63
34:SR:238:ASP:N	34:SR:238:ASP:OD1	2.32	0.63
58:N2:94:ARG:NH2	36:5:1757:A:OP1	128.24	0.63
16:C4:11:SER:OG	16:C4:12:GLN:N	4.14	0.63
24:D2:56:HIS:O	1:6:861:U:O2'	356.30	0.63
28:D6:58:VAL:HG22	28:D6:59:TYR:H	4.10	0.63
29:D7:53:ALA:HB1	29:D7:62:ILE:HD11	3.29	0.63
30:D8:29:ARG:HA	30:D8:41:VAL:HA	1.81	0.63
43:L6:56:LYS:HD2	43:L6:98:VAL:HG13	1.81	0.63
44:L7:64:GLN:HA	44:L7:67:ARG:HD2	2.75	0.63
43:L6:3:ALA:HB2	68:O2:77:ALA:HB2	1.81	0.63
74:O8:22:THR:HG22	74:O8:74:LYS:HB3	4.11	0.63
39:L2:108:PRO:HG2	79:Q3:86:LEU:HD22	1.81	0.63
4:S2:89:GLN:HG3	4:S2:93:GLY:HA2	6.31	0.63
34:SR:76:ASP:OD1	34:SR:76:ASP:N	2.28	0.63
1:2:1738:U:H2'	1:2:1739:C:C6	2.34	0.62
1:2:720:G:H1'	1:2:721:U:H5''	1.80	0.62
21:C9:63:ARG:NH1	1:6:1481:C:OP2	404.08	0.62
28:D6:79:ILE:HG23	28:D6:84:VAL:HG21	1.80	0.62
37:3:121:U:C2	42:L5:268:GLU:HB3	2.34	0.62
51:M5:53:TYR:HD1	51:M5:133:ILE:HD13	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
74:O8:26:LYS:HE2	36:5:1751:G:H5''	127.78	0.62
6:S4:125:LYS:NZ	6:S4:225:VAL:O	2.25	0.62
36:1:148:G:OP2	51:M5:4:TYR:OH	2.12	0.62
36:1:2724:U:OP1	57:N1:57:TYR:OH	2.10	0.62
1:6:1537:C:O2'	1:6:1540:G:O6	2.17	0.62
17:C5:79:HIS:O	17:C5:81:ARG:N	2.32	0.62
28:D6:23:CYS:CB	28:D6:74:CYS:SG	4.58	0.62
39:L2:227:ARG:HG2	39:L2:239:ALA:HB2	1.81	0.62
42:L5:146:LEU:HB3	36:5:2746:A:H2	258.69	0.62
55:M9:68:GLN:NE2	55:M9:72:GLU:OE2	4.43	0.62
63:N7:55:LYS:O	63:N7:57:HIS:N	3.21	0.62
67:O1:13:THR:HG23	67:O1:72:ARG:HH11	1.65	0.62
1:2:858:G:OP1	9:S7:116:ARG:NH2	2.31	0.62
15:C3:73:ARG:O	15:C3:77:SER:OG	2.11	0.62
17:C5:115:TYR:N	17:C5:118:GLU:OE1	2.32	0.62
26:D4:55:VAL:HG12	26:D4:75:VAL:HG22	7.21	0.62
40:L3:25:ILE:H	40:L3:25:ILE:HD13	1.64	0.62
41:L4:38:VAL:HG13	41:L4:113:VAL:HG11	2.03	0.62
42:L5:218:ARG:HH21	42:L5:222:LEU:HD11	1.64	0.62
53:M7:26:PHE:HE1	53:M7:120:ASN:HA	1.65	0.62
63:N7:23:VAL:HG12	63:N7:45:GLY:HA3	1.81	0.62
72:O6:63:ASN:O	72:O6:65:GLY:N	4.70	0.62
73:O7:21:ARG:HD2	73:O7:39:TYR:HB2	1.80	0.62
74:O8:16:ARG:HH11	74:O8:70:PRO:HG3	3.65	0.62
36:1:1615:C:H2'	36:1:1616:U:C6	2.35	0.62
39:L2:201:GLY:O	39:L2:204:MET:HG2	1.98	0.62
58:N2:50:LEU:O	58:N2:52:ASN:N	2.32	0.62
45:L8:26:LEU:HD21	63:N7:123:GLN:HG2	1.80	0.62
6:S4:211:LYS:NZ	6:S4:212:ASP:O	2.32	0.62
7:S5:91:GLU:O	7:S5:95:ASN:ND2	2.32	0.62
34:SR:258:THR:O	34:SR:275:ARG:NH1	2.32	0.62
76:Q0:111:ARG:NH2	36:5:3121:U:OP2	321.12	0.62
15:C3:5:HIS:HB3	15:C3:117:LEU:HD13	1.80	0.62
19:C7:5:ARG:O	19:C7:10:LYS:NZ	4.19	0.62
20:C8:5:VAL:HG23	27:D5:42:LEU:HD21	1.81	0.62
41:L4:251:THR:O	41:L4:253:ALA:N	3.29	0.62
50:M4:73:PRO:HG2	50:M4:76:ALA:HB2	3.06	0.62
8:S6:98:ARG:HD3	8:S6:99:GLY:N	2.28	0.62
11:S9:65:LYS:HG2	11:S9:70:LEU:HD21	1.81	0.62
11:S9:72:GLU:OE2	1:6:761:G:O2'	397.07	0.62
35:SM:104:LYS:O	35:SM:108:GLN:HG2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1591:G:O2'	36:1:1799:A:N1	2.30	0.62
38:4:57:C:OP2	73:O7:68:LYS:NZ	2.33	0.62
68:O2:27:ARG:NH1	36:5:1433:A:N3	169.56	0.62
40:L3:126:LYS:NZ	36:5:3294:A:OP2	189.31	0.62
49:M3:58:VAL:HG13	36:5:75:G:H5''	88.32	0.62
1:6:428:A:N3	1:6:440:U:O2'	2.25	0.62
15:C3:64:ARG:NH2	1:6:861:U:OP1	347.10	0.62
30:D8:19:THR:HB	30:D8:66:LEU:HB2	1.80	0.62
46:L9:149:ASN:N	46:L9:149:ASN:OD1	2.32	0.62
36:1:686:G:OP2	49:M3:39:ARG:NH2	2.33	0.62
58:N2:14:THR:HG23	58:N2:66:VAL:HG22	3.70	0.62
59:N3:17:LEU:HD11	59:N3:98:ASN:HB3	1.82	0.62
69:O3:75:HIS:HB3	69:O3:80:VAL:HG12	1.81	0.62
69:O3:9:VAL:HG21	69:O3:44:TYR:HE1	2.09	0.62
8:S6:10:ASN:HB3	8:S6:128:THR:HA	2.66	0.62
1:2:79:C:H1'	8:S6:174:LYS:HD3	1.81	0.62
36:1:1925:U:O2'	36:1:1927:G:N7	2.32	0.62
36:1:2216:G:OP1	72:O6:75:LYS:NZ	2.30	0.62
36:1:3122:A:N1	46:L9:70:THR:HG21	2.15	0.62
1:2:584:C:H1'	32:E0:18:THR:HG21	1.80	0.62
1:6:1280:C:H2'	1:6:1281:G:C8	2.35	0.62
22:D0:63:LEU:HD12	22:D0:84:MET:HB3	4.69	0.62
2:S0:52:LYS:HB3	23:D1:82:VAL:HG22	1.81	0.62
36:1:3243:A:N6	52:M6:157:GLU:OE2	2.24	0.62
56:N0:91:TYR:O	56:N0:137:ARG:NH1	2.81	0.62
67:O1:26:LYS:HE2	67:O1:64:VAL:HG21	1.82	0.62
75:O9:21:ARG:HD3	75:O9:22:PRO:O	2.00	0.62
2:S0:26:ALA:HB3	2:S0:149:LEU:HB2	2.45	0.62
5:S3:115:ILE:O	5:S3:119:ALA:N	2.95	0.62
36:1:1256:G:O6	36:1:1261:G:N2	2.32	0.62
36:1:1480:G:H4'	36:1:1481:A:OP1	2.00	0.62
36:1:1740:U:H1'	36:1:1741:A:H2	1.64	0.62
36:1:2149:A:N6	36:1:2187:G:O2'	2.33	0.62
1:2:772:G:N2	1:2:774:A:O2'	2.32	0.62
1:6:1561:U:H4'	1:6:1599:C:H4'	1.81	0.62
12:C0:11:ILE:HD13	12:C0:35:ILE:HG21	1.82	0.62
21:C9:86:ARG:NH1	21:C9:90:PRO:O	2.32	0.62
41:L4:234:ASN:OD1	41:L4:235:LEU:N	2.32	0.62
47:M0:177:ASP:OD2	47:M0:177:ASP:N	3.76	0.62
47:M0:84:ALA:O	47:M0:140:THR:HG22	2.00	0.62
49:M3:74:GLY:O	49:M3:101:ARG:NH1	2.44	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:M8:165:ILE:HD13	54:M8:167:SER:H	6.55	0.62
55:M9:28:GLU:HG3	55:M9:49:THR:HG22	5.13	0.62
62:N6:3:LYS:NZ	62:N6:8:VAL:O	2.26	0.62
74:O8:16:ARG:NH1	74:O8:70:PRO:HG3	4.42	0.62
10:S8:166:TYR:HB3	10:S8:184:LEU:HD22	1.81	0.62
34:SR:31:ASN:N	34:SR:31:ASN:OD1	2.32	0.62
36:1:1103:A:O2'	36:1:1104:G:OP1	2.15	0.62
36:1:585:A:H5'	69:O3:70:LYS:HE2	1.82	0.62
1:2:100:A:H61	1:2:385:A:H1'	1.65	0.62
1:2:207:U:O2	10:S8:178:ARG:NH1	2.32	0.62
36:5:1781:C:H2'	36:5:1782:U:C6	2.35	0.62
19:C7:21:TYR:OH	19:C7:62:GLN:OE1	2.23	0.62
19:C7:6:THR:OG1	19:C7:7:LYS:N	2.47	0.62
2:S0:88:LYS:NZ	19:C7:82:ASP:OD1	3.85	0.62
2:S0:88:LYS:HE2	2:S0:201:LEU:HG	6.02	0.62
6:S4:19:LEU:HD11	6:S4:108:ARG:HD2	1.81	0.62
7:S5:50:GLU:O	7:S5:65:ARG:NH2	2.33	0.62
36:1:2185:G:O2'	36:1:2314:U:OP2	2.17	0.62
36:1:2836:C:H5	36:1:2852:C:H42	1.48	0.62
36:5:1783:U:H2'	36:5:1784:G:H8	1.64	0.62
36:5:2267:C:H2'	36:5:2268:U:H6	1.64	0.62
1:6:1398:U:H3'	1:6:1399:C:H4'	1.82	0.62
1:6:1603:U:H2'	1:6:1604:U:C6	2.35	0.62
18:C6:120:ASP:OD1	18:C6:121:SER:N	2.33	0.62
36:1:1794:G:H4'	39:L2:191:LEU:HD13	1.80	0.62
44:L7:60:ARG:HA	44:L7:63:ILE:HD12	2.41	0.62
45:L8:244:ALA:HA	45:L8:247:ASP:HB2	2.32	0.62
46:L9:22:SER:OG	46:L9:39:LYS:NZ	2.32	0.62
50:M4:22:LEU:HB3	50:M4:64:VAL:HG13	6.00	0.62
41:L4:106:TRP:HB2	51:M5:199:LEU:HD12	1.81	0.62
54:M8:64:VAL:HG11	54:M8:113:LYS:HD3	2.66	0.62
62:N6:32:SER:HA	62:N6:49:PRO:HA	2.69	0.62
72:O6:15:LYS:HA	72:O6:15:LYS:HZ3	5.39	0.62
78:Q2:12:CYS:SG	78:Q2:17:CYS:CB	2.68	0.62
79:Q3:38:ASP:HA	79:Q3:45:LYS:HA	1.82	0.62
1:2:399:A:N3	6:S4:3:ARG:NH1	2.48	0.62
10:S8:8:ARG:NH2	10:S8:21:PHE:H	1.98	0.62
34:SR:216:LYS:HA	34:SR:239:GLU:HG3	2.14	0.62
36:1:999:G:N3	36:1:1002:A:N6	2.48	0.61
36:5:2403:G:H5'	36:5:2872:A:N1	2.15	0.61
40:L3:62:ARG:NH1	36:5:3039:C:OP1	275.70	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1054:U:H2'	1:6:1055:U:H6	1.64	0.61
17:C5:126:VAL:HB	1:6:1459:C:H5'	344.15	0.61
21:C9:105:LEU:HD13	21:C9:122:ARG:HD3	1.99	0.61
40:L3:296:THR:OG1	40:L3:357:LYS:O	2.18	0.61
40:L3:41:VAL:HG22	40:L3:185:GLY:HA3	1.81	0.61
46:L9:70:THR:HG21	36:5:3122:A:N1	323.98	0.61
35:SM:59:GLY:O	35:SM:63:ASP:N	2.84	0.61
36:1:2115:G:H22	36:1:2120:A:H1'	1.63	0.61
36:1:3041:U:OP1	59:N3:12:ARG:NH1	2.32	0.61
1:2:483:A:H61	1:2:505:A:H1'	1.65	0.61
1:2:904:G:H2'	1:2:905:A:H8	1.65	0.61
36:5:2219:A:H2'	36:5:2220:A:C8	2.35	0.61
17:C5:40:ARG:NH1	1:6:1556:A:O2'	385.31	0.61
21:C9:37:VAL:HG11	21:C9:100:ILE:HD11	1.82	0.61
39:L2:193:ARG:NH2	36:5:2181:C:OP1	197.92	0.61
40:L3:227:GLU:HG3	40:L3:270:ARG:HE	4.17	0.61
46:L9:62:ARG:O	46:L9:66:ALA:N	2.78	0.61
58:N2:104:ARG:HH11	58:N2:106:ALA:HB2	3.35	0.61
67:O1:74:ARG:HH12	67:O1:109:VAL:HG11	5.38	0.61
11:S9:92:LYS:HB2	11:S9:95:TYR:HD2	6.52	0.61
36:1:2592:G:H4'	36:1:2594:C:C2	2.34	0.61
1:2:75:U:H3'	1:2:76:A:O4'	2.00	0.61
1:2:789:A:O2'	6:S4:106:LYS:NZ	2.31	0.61
36:5:1757:A:H2'	36:5:1758:G:H8	1.66	0.61
1:6:1226:A:O2'	1:6:1256:A:N6	2.33	0.61
1:6:500:C:O2'	1:6:501:U:O4'	2.19	0.61
1:6:58:U:O2'	1:6:451:A:N3	2.31	0.61
18:C6:50:GLU:OE1	18:C6:114:ARG:NH1	2.33	0.61
19:C7:26:LEU:O	19:C7:59:LYS:NZ	6.25	0.61
1:2:1616:G:O2'	30:D8:18:ARG:NH1	2.32	0.61
46:L9:48:VAL:HG13	46:L9:52:LEU:HB3	1.81	0.61
49:M3:175:SER:O	49:M3:178:LYS:N	2.33	0.61
52:M6:133:ARG:NE	36:5:1316:C:OP2	293.69	0.61
55:M9:105:LEU:HD21	55:M9:139:VAL:HG13	6.78	0.61
2:S0:120:LEU:HD21	2:S0:144:ILE:HD11	3.00	0.61
17:C5:130:ARG:HH22	35:SM:70:ASN:HB3	3.98	0.61
36:1:2155:G:O2'	39:L2:227:ARG:NH2	2.32	0.61
36:1:2986:U:H2'	36:1:2987:A:H8	1.64	0.61
1:2:38:C:H2'	1:2:39:A:H5'	1.80	0.61
1:2:734:A:H5''	1:2:735:C:OP1	2.01	0.61
76:Q0:111:ARG:NH2	36:5:3120:C:H3'	321.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1380:U:O2'	1:6:1516:A:N1	2.33	0.61
38:8:79:A:H3'	38:8:80:A:C8	2.36	0.61
39:L2:172:GLY:HA3	79:Q3:68:ALA:N	4.73	0.61
41:L4:351:PRO:HA	44:L7:71:ALA:HA	1.82	0.61
42:L5:156:GLY:HA2	42:L5:181:PRO:HD3	1.82	0.61
44:L7:125:GLU:OE2	44:L7:128:LYS:NZ	2.28	0.61
50:M4:23:ILE:HA	50:M4:63:VAL:HG23	1.83	0.61
68:O2:123:LYS:HA	68:O2:126:LEU:HD12	1.82	0.61
71:O5:119:LYS:NZ	71:O5:119:LYS:HA	4.89	0.61
72:O6:2:THR:OG1	72:O6:3:VAL:N	2.46	0.61
3:S1:212:VAL:O	3:S1:214:LYS:N	2.31	0.61
4:S2:55:GLU:O	4:S2:59:HIS:ND1	2.33	0.61
8:S6:120:GLU:HG3	8:S6:125:THR:HB	2.25	0.61
8:S6:164:LYS:N	8:S6:167:LYS:O	2.31	0.61
36:1:787:G:H2'	36:1:788:C:C6	2.35	0.61
1:2:520:A:H2'	1:2:521:A:C8	2.35	0.61
36:5:2254:U:H2'	36:5:2261:G:N2	2.16	0.61
16:C4:18:ARG:NH2	16:C4:31:THR:OG1	6.87	0.61
20:C8:4:VAL:HG11	27:D5:82:HIS:HB2	5.53	0.61
40:L3:173:GLN:O	40:L3:175:LYS:N	2.34	0.61
40:L3:282:ILE:HD13	40:L3:322:ILE:HD12	1.83	0.61
41:L4:304:GLN:O	41:L4:306:THR:N	2.33	0.61
53:M7:25:SER:HB3	53:M7:28:ASN:HB2	1.95	0.61
72:O6:34:SER:HG	72:O6:37:THR:HG1	1.47	0.61
74:O8:65:LEU:HD23	74:O8:68:SER:HB2	1.82	0.61
36:5:1366:A:C2	36:5:1367:G:C4	2.89	0.61
1:6:698:U:H3	1:6:740:A:H61	1.47	0.61
1:6:222:A:H62	1:6:833:U:H3	1.48	0.61
18:C6:37:THR:O	18:C6:45:ARG:NH1	2.46	0.61
22:D0:39:SER:HA	22:D0:42:VAL:HG12	1.83	0.61
24:D2:104:LEU:HB2	24:D2:125:ILE:HA	1.82	0.61
33:E1:108:VAL:HB	33:E1:114:VAL:HG22	1.83	0.61
1:2:1236:A:H1'	33:E1:138:ARG:HH12	1.63	0.61
49:M3:166:ALA:HB1	64:N8:147:LEU:HD21	2.16	0.61
67:O1:62:ARG:HB2	67:O1:66:GLY:O	2.28	0.61
4:S2:41:LEU:HD12	4:S2:68:ILE:HD13	1.82	0.61
5:S3:222:VAL:HG11	34:SR:229:LYS:HA	3.85	0.61
9:S7:96:ARG:NH2	9:S7:128:ASP:OD2	2.34	0.61
34:SR:165:ASP:O	34:SR:184:ASN:ND2	2.34	0.61
36:1:129:U:H2'	36:1:130:A:C8	2.35	0.61
36:1:796:U:H2'	36:1:797:U:H6	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:304:U:H2'	1:2:305:C:C6	2.34	0.61
1:2:514:G:H1	1:2:543:C:N4	1.97	0.61
1:6:44:U:OP2	1:6:437:A:N6	2.33	0.61
1:6:792:U:H3'	1:6:793:A:C8	2.36	0.61
2:S0:7:PHE:CZ	23:D1:43:GLY:HA2	3.11	0.61
41:L4:11:LEU:HD13	41:L4:159:ILE:HD11	1.82	0.61
37:3:121:U:N3	42:L5:268:GLU:HB3	2.15	0.61
47:M0:191:LYS:HB3	47:M0:213:PHE:CE2	2.33	0.61
55:M9:41:ILE:HG23	55:M9:50:ILE:HD12	1.82	0.61
60:N4:23:ARG:NH2	60:N4:25:ASP:OD1	3.34	0.61
62:N6:45:ILE:HD12	62:N6:119:ILE:HG23	1.81	0.61
63:N7:36:HIS:O	63:N7:38:PHE:N	2.29	0.61
39:L2:172:GLY:HA3	79:Q3:68:ALA:H	3.98	0.61
2:S0:148:ASP:OD1	2:S0:149:LEU:N	2.40	0.61
36:1:1095:U:H4'	36:1:1096:U:H5''	1.83	0.61
1:2:1511:U:H2'	1:2:1512:G:C8	2.35	0.61
1:2:1551:U:H3'	17:C5:43:ARG:HH22	1.66	0.61
55:M9:20:ARG:HD2	36:5:1874:A:OP2	141.93	0.61
1:6:1045:C:H2'	1:6:1046:G:C8	2.36	0.61
1:6:621:A:N3	1:6:1107:G:H1'	2.14	0.61
18:C6:18:ALA:HB2	18:C6:69:VAL:HG13	1.97	0.61
23:D1:71:ARG:HG3	23:D1:83:TRP:CZ2	3.48	0.61
25:D3:69:ARG:NH1	25:D3:116:ASP:OD2	2.93	0.61
1:2:930:A:H4'	28:D6:70:LYS:HG3	1.83	0.61
1:2:545:A:H2'	32:E0:31:LYS:HD2	1.81	0.61
58:N2:48:GLY:O	58:N2:50:LEU:N	3.64	0.61
66:O0:9:SER:OG	66:O0:10:ILE:N	3.39	0.61
36:1:1926:C:H2'	79:Q3:7:LYS:HE2	1.81	0.61
5:S3:34:TYR:HE2	5:S3:37:VAL:HG13	2.25	0.61
1:2:755:A:H2	6:S4:13:ALA:HA	1.64	0.61
11:S9:92:LYS:O	11:S9:94:ASP:N	2.30	0.61
1:2:121:U:H1'	6:S4:33:ALA:HB3	1.83	0.61
1:2:1583:A:OP1	18:C6:135:ARG:NH2	2.27	0.61
36:5:2268:U:H2'	36:5:2269:U:H2'	1.82	0.61
36:5:2448:G:H1	36:5:2498:U:H3	1.49	0.61
1:6:1218:G:H5''	1:6:1444:A:H61	1.66	0.61
12:C0:38:LYS:HD3	12:C0:41:TYR:CZ	2.36	0.61
18:C6:40:GLU:HA	18:C6:42:GLU:N	2.16	0.61
28:D6:87:ARG:NH2	28:D6:94:ASN:O	2.34	0.61
42:L5:187:THR:HG22	42:L5:189:GLU:H	1.65	0.61
44:L7:88:ARG:NH1	44:L7:91:GLY:O	2.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:M7:59:PRO:HG3	53:M7:76:PHE:CD1	2.35	0.61
36:1:2556:C:O2'	63:N7:135:ARG:NH2	2.33	0.61
2:S0:103:THR:O	2:S0:106:SER:OG	2.18	0.61
34:SR:95:ALA:O	34:SR:97:GLY:N	4.83	0.61
36:1:651:G:O2'	36:1:1435:A:OP1	2.19	0.61
36:1:2676:A:N1	48:M1:22:SER:OG	2.29	0.61
1:2:1000:C:N4	1:2:1003:A:OP2	2.32	0.61
36:5:2102:U:H2'	36:5:2103:U:C6	2.36	0.61
36:5:3287:U:H2'	36:5:3288:G:H5'	1.83	0.61
1:2:917:U:O2	16:C4:41:ARG:NH1	2.34	0.61
24:D2:82:LYS:O	24:D2:84:GLY:N	2.31	0.61
27:D5:71:ILE:HD12	27:D5:76:ALA:HA	1.83	0.61
14:C2:54:ARG:HH21	33:E1:127:GLY:H	1.46	0.61
42:L5:153:THR:HG22	42:L5:179:ARG:HH11	1.65	0.61
52:M6:183:ALA:HA	52:M6:186:ALA:HB2	2.51	0.61
68:O2:21:HIS:CE1	68:O2:24:ARG:HD2	2.51	0.61
75:O9:23:LEU:HD22	75:O9:24:PRO:HD2	1.96	0.61
79:Q3:87:ARG:O	79:Q3:91:GLU:HG2	3.80	0.61
2:S0:74:VAL:HG23	2:S0:118:PRO:HB3	2.00	0.61
5:S3:113:LEU:HD12	5:S3:117:ARG:HD3	5.11	0.61
9:S7:50:ASP:HA	9:S7:56:LYS:HG2	3.40	0.61
9:S7:74:GLN:HE22	9:S7:92:PHE:HB2	2.10	0.61
1:2:1230:A:H2'	1:2:1258:U:H5	1.66	0.60
36:5:2400:G:O2'	36:5:2401:A:OP1	2.18	0.60
7:S5:102:ARG:NH1	1:6:1473:U:O2'	354.04	0.60
14:C2:87:PRO:HA	14:C2:140:PHE:HE1	1.66	0.60
17:C5:37:ALA:O	17:C5:42:ARG:NH1	2.70	0.60
18:C6:107:LYS:O	18:C6:111:SER:OG	4.14	0.60
20:C8:28:ILE:HD12	20:C8:61:LEU:HD11	1.83	0.60
49:M3:59:ARG:HD2	49:M3:69:VAL:HG23	4.80	0.60
59:N3:2:SER:HA	59:N3:56:ASP:HA	5.18	0.60
36:1:2818:U:O5'	65:N9:2:ALA:HB2	2.01	0.60
3:S1:107:THR:HB	3:S1:111:ARG:HH21	1.66	0.60
3:S1:22:ASP:O	3:S1:25:THR:OG1	4.54	0.60
36:1:58:G:H4'	51:M5:155:VAL:HG12	1.83	0.60
36:1:743:C:N3	54:M8:141:ARG:NH1	2.49	0.60
1:2:1553:G:N2	1:2:1555:A:H3'	2.16	0.60
59:N3:92:PHE:CE1	36:5:3051:U:H1'	245.32	0.60
14:C2:68:GLU:O	14:C2:70:ASN:ND2	2.35	0.60
28:D6:10:ARG:NE	1:6:1795:U:O2	328.46	0.60
29:D7:61:THR:HG23	29:D7:62:ILE:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:E1:109:ASP:O	33:E1:111:GLU:N	2.85	0.60
3:S1:165:ARG:O	3:S1:169:SER:OG	2.16	0.60
34:SR:63:GLY:HA3	34:SR:92:TRP:HH2	2.59	0.60
36:1:1345:G:N2	41:L4:307:GLN:OE1	2.33	0.60
36:1:3355:U:O2'	36:1:3357:U:OP2	2.18	0.60
1:2:407:A:OP1	8:S6:94:ARG:NE	2.33	0.60
56:N0:137:ARG:NH1	36:5:1213:G:OP1	325.03	0.60
36:5:1952:G:H1	36:5:2094:C:N4	1.98	0.60
42:L5:270:LYS:HD3	37:7:2:G:H4'	321.31	0.60
42:L5:4:GLN:N	42:L5:4:GLN:OE1	2.29	0.60
55:M9:39:ASN:HA	55:M9:42:ARG:HH11	4.41	0.60
59:N3:87:ARG:NE	59:N3:121:GLU:OE2	2.34	0.60
36:1:1603:A:H61	61:N5:71:THR:HG21	1.66	0.60
67:O1:31:ARG:HH11	67:O1:31:ARG:HB3	4.21	0.60
72:O6:66:GLU:OE2	72:O6:91:ASN:ND2	3.08	0.60
73:O7:64:MET:HB3	73:O7:68:LYS:HB3	4.27	0.60
2:S0:13:ASP:OD1	2:S0:179:ARG:NH2	4.07	0.60
36:1:1630:U:OP1	63:N7:67:LYS:NZ	2.29	0.60
36:1:2617:U:H5	36:1:2621:G:OP2	1.83	0.60
60:N4:34:SER:OG	36:5:3085:G:OP1	229.11	0.60
36:5:67:A:O2'	36:5:315:C:O2	2.18	0.60
49:M3:32:LYS:NZ	36:5:686:G:N7	85.19	0.60
38:8:125:U:O2'	38:8:126:A:H5'	2.02	0.60
38:8:157:U:H2'	38:8:158:U:H6	1.66	0.60
71:O5:7:TYR:CE2	38:8:86:U:H2'	20.00	0.60
15:C3:22:ALA:HB1	15:C3:23:PRO:HA	2.06	0.60
41:L4:38:VAL:HG21	41:L4:121:ALA:HB2	2.68	0.60
41:L4:282:SER:HB3	54:M8:126:GLN:HE21	1.74	0.60
4:S2:77:GLN:H	4:S2:77:GLN:HE21	1.49	0.60
8:S6:153:VAL:O	8:S6:155:ASP:N	3.41	0.60
9:S7:35:LYS:O	9:S7:37:GLU:N	2.34	0.60
10:S8:56:ARG:HH22	1:6:332:U:P	286.85	0.60
1:2:1426:C:H5''	35:SM:93:ARG:NH1	2.16	0.60
36:1:1308:A:H8	36:1:1308:A:OP2	1.85	0.60
1:2:751:G:H2'	1:2:752:A:C8	2.36	0.60
36:5:2207:A:H2'	36:5:2208:A:O4'	2.01	0.60
12:C0:14:TYR:OH	12:C0:34:GLU:OE1	2.13	0.60
40:L3:141:GLY:O	40:L3:143:GLY:N	2.35	0.60
41:L4:186:LYS:N	41:L4:200:THR:O	2.70	0.60
60:N4:63:ILE:O	60:N4:65:GLU:N	4.91	0.60
61:N5:82:LEU:HD11	61:N5:135:ILE:HG21	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
69:O3:45:LEU:HD11	69:O3:73:ARG:HA	3.01	0.60
71:O5:31:LEU:O	71:O5:35:LYS:N	3.13	0.60
74:O8:12:LEU:O	74:O8:15:THR:OG1	2.54	0.60
74:O8:42:LYS:HG2	74:O8:55:VAL:HG22	1.82	0.60
3:S1:82:ARG:NH1	3:S1:191:GLU:OE2	4.21	0.60
4:S2:152:HIS:ND1	4:S2:174:ARG:HG2	2.88	0.60
6:S4:73:ASP:HB3	6:S4:164:LEU:HD22	1.84	0.60
7:S5:135:ASP:O	7:S5:139:ASN:ND2	2.31	0.60
7:S5:40:ILE:HG23	7:S5:42:LEU:HD22	1.81	0.60
9:S7:133:THR:OG1	9:S7:134:GLU:N	2.32	0.60
36:1:1286:A:O2'	36:1:1287:A:OP2	2.18	0.60
36:1:1820:U:O2	36:1:1822:C:N4	2.34	0.60
36:1:241:G:O6	36:1:242:C:N4	2.35	0.60
1:2:154:G:OP1	8:S6:2:LYS:NZ	2.34	0.60
37:3:97:A:H2'	37:3:98:C:H6	1.65	0.60
1:6:138:A:OP2	1:6:1706:C:O2'	2.18	0.60
1:6:1474:G:H2'	1:6:1475:A:C8	2.36	0.60
1:6:885:G:H2'	1:6:886:U:C6	2.37	0.60
37:7:112:G:H2'	37:7:113:C:C6	2.35	0.60
23:D1:5:LYS:HD3	23:D1:5:LYS:H	1.67	0.60
25:D3:102:VAL:HG12	25:D3:127:VAL:HB	1.83	0.60
41:L4:158:SER:HA	41:L4:213:ASN:HB2	1.83	0.60
42:L5:148:ILE:HG23	42:L5:151:GLN:HB3	1.93	0.60
47:M0:49:CYS:HB2	47:M0:172:GLY:HA2	1.83	0.60
48:M1:28:ASP:HB3	48:M1:32:ARG:HH21	1.67	0.60
56:N0:155:ARG:HD3	56:N0:172:TYR:CG	5.39	0.60
56:N0:77:VAL:HG11	56:N0:106:LEU:HG	2.87	0.60
66:O0:78:GLY:HA2	66:O0:87:VAL:HG13	2.19	0.60
72:O6:59:ASP:O	72:O6:63:ASN:ND2	4.98	0.60
36:1:1841:A:N3	75:O9:45:ARG:NH2	2.48	0.60
7:S5:140:THR:HA	7:S5:214:LYS:HD2	2.06	0.60
10:S8:152:ILE:H	10:S8:152:ILE:HD13	4.51	0.60
36:1:2572:C:O2'	36:1:2573:G:O4'	2.19	0.60
36:5:1596:C:H2'	36:5:1597:C:C6	2.37	0.60
54:M8:21:SER:OG	36:5:673:U:OP1	150.51	0.60
3:S1:27:LYS:NZ	1:6:896:U:OP1	254.50	0.60
40:L3:43:LEU:HG	40:L3:181:ILE:HG21	2.13	0.60
41:L4:341:SER:O	36:5:515:C:O2'	298.65	0.60
44:L7:121:LYS:HB2	57:N1:133:ALA:HB3	1.85	0.60
36:1:1841:A:H1'	75:O9:45:ARG:HH12	1.66	0.60
78:Q2:17:CYS:HB2	78:Q2:77:CYS:SG	2.40	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S2:69:ILE:HG12	4:S2:133:LYS:HB3	2.47	0.60
1:2:121:U:O2'	6:S4:33:ALA:O	2.16	0.60
9:S7:174:ASN:ND2	9:S7:180:GLN:OE1	2.30	0.60
1:2:74:U:H1'	1:2:75:U:H4'	1.83	0.60
47:M0:150:GLU:O	47:M0:152:LEU:N	3.98	0.60
52:M6:54:TYR:HE2	52:M6:58:LEU:HD13	2.22	0.60
54:M8:165:ILE:HD13	54:M8:166:LEU:H	5.53	0.60
61:N5:46:TYR:HB3	71:O5:75:TYR:HB3	1.82	0.60
2:S0:167:LYS:HG2	2:S0:168:HIS:CD2	3.33	0.60
5:S3:31:GLU:O	5:S3:54:ARG:NH2	3.00	0.60
10:S8:159:GLN:OE1	10:S8:166:TYR:N	2.35	0.60
36:1:114:A:N1	36:1:266:A:O2'	2.32	0.60
36:1:2163:C:H4'	39:L2:7:ASN:O	2.02	0.60
36:1:2703:A:OP2	42:L5:23:ARG:NH1	2.34	0.60
1:2:1615:C:O2'	1:2:1616:G:OP2	2.19	0.60
1:2:968:U:O3'	1:2:1032:G:N2	2.35	0.60
36:5:1012:G:N2	36:5:1038:C:N3	2.46	0.60
36:5:1152:G:N2	36:5:1200:A:H61	2.00	0.60
36:5:1818:U:H2'	36:5:1819:U:O4'	2.02	0.60
76:Q0:111:ARG:HH21	36:5:3120:C:H3'	321.50	0.60
19:C7:7:LYS:N	1:6:1316:G:OP1	409.99	0.60
1:6:1698:G:O2'	1:6:1699:G:O5'	2.19	0.60
18:C6:14:LYS:O	18:C6:123:ARG:NH1	2.28	0.60
41:L4:193:LYS:HA	41:L4:198:ARG:HA	1.84	0.60
5:S3:69:LEU:HA	5:S3:72:LEU:HD12	2.33	0.60
8:S6:148:SER:HB3	60:N4:98:PRO:HG3	1.82	0.60
34:SR:211:ILE:HD11	34:SR:225:LEU:HD13	1.84	0.60
36:1:2747:A:OP1	42:L5:176:SER:OG	2.18	0.60
36:5:999:G:N3	36:5:1002:A:N6	2.50	0.60
36:5:2105:G:H2'	36:5:2106:A:H8	1.65	0.60
25:D3:97:ASP:N	25:D3:100:ASP:OD2	4.25	0.60
25:D3:56:LYS:HG3	25:D3:98:GLU:HG3	1.84	0.60
42:L5:110:LEU:HA	42:L5:113:LEU:HB2	3.08	0.60
36:1:1677:G:N7	58:N2:74:LYS:NZ	2.50	0.60
45:L8:165:PHE:HA	72:O6:47:ILE:HD13	2.47	0.60
9:S7:105:THR:O	9:S7:107:ARG:N	4.08	0.60
9:S7:55:LYS:HE2	9:S7:87:ASP:HA	2.60	0.60
36:1:305:U:C5	36:1:2776:C:H1'	2.37	0.59
36:1:2927:C:H2'	36:1:2928:C:C6	2.36	0.59
1:2:319:U:H5'	1:2:320:U:H5	1.66	0.59
36:5:161:G:N2	36:5:260:C:O2	2.26	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:O7:55:ARG:NH1	36:5:353:G:N7	110.51	0.59
17:C5:64:LYS:HA	17:C5:73:PRO:HB3	1.83	0.59
20:C8:18:LEU:HD12	20:C8:102:ALA:HB2	2.40	0.59
21:C9:10:ALA:HB3	21:C9:13:ASP:HB2	2.04	0.59
25:D3:35:GLY:O	25:D3:37:ALA:N	2.35	0.59
28:D6:85:ARG:O	1:6:1797:A:N6	341.36	0.59
33:E1:121:CYS:HB3	33:E1:130:VAL:HG11	6.03	0.59
42:L5:69:ILE:HD12	42:L5:69:ILE:H	4.96	0.59
47:M0:9:TYR:O	47:M0:59:GLN:NE2	2.35	0.59
63:N7:102:GLU:H	63:N7:107:ARG:HH21	1.50	0.59
10:S8:114:GLU:HG2	10:S8:120:THR:HA	1.84	0.59
34:SR:116:ASP:HA	34:SR:156:VAL:HG11	1.84	0.59
36:1:1064:A:H4'	36:1:1065:A:O5'	2.01	0.59
36:1:2107:A:H2	36:1:3344:A:H8	1.50	0.59
36:1:975:C:H2'	36:1:976:U:C6	2.37	0.59
72:O6:25:LYS:HB3	36:5:156:G:OP2	88.10	0.59
78:Q2:10:THR:OG1	36:5:2714:G:OP2	218.51	0.59
22:D0:74:GLU:HG2	1:6:1429:G:H1'	377.49	0.59
1:2:522:U:O2'	26:D4:60:PHE:O	2.16	0.59
42:L5:85:ARG:NH2	42:L5:250:ASP:OD1	2.32	0.59
44:L7:102:VAL:HA	44:L7:105:LEU:HD12	1.84	0.59
56:N0:1:MET:HE1	56:N0:32:SER:N	2.17	0.59
66:O0:95:ALA:HB2	66:O0:101:LEU:HD23	1.84	0.59
67:O1:83:GLU:O	67:O1:85:ALA:N	3.77	0.59
62:N6:127:GLU:O	71:O5:68:GLN:NE2	49.07	0.59
75:O9:21:ARG:NH1	75:O9:22:PRO:O	2.30	0.59
6:S4:252:ARG:HB2	11:S9:71:PHE:HE2	1.66	0.59
7:S5:166:ARG:NH1	7:S5:170:GLN:OE1	2.35	0.59
11:S9:110:GLN:NE2	11:S9:126:ARG:HG2	2.17	0.59
11:S9:38:ASN:HB2	11:S9:40:LYS:H	1.67	0.59
35:SM:48:ARG:HH22	36:5:1017:C:H4'	337.12	0.59
34:SR:16:HIS:CE1	34:SR:43:ILE:HG12	2.37	0.59
36:1:551:A:O2'	36:1:552:G:O5'	2.19	0.59
1:6:1120:U:H2'	1:6:1121:C:C6	2.36	0.59
13:C1:111:VAL:HG23	13:C1:139:VAL:HG21	3.27	0.59
20:C8:99:HIS:CD2	20:C8:101:LEU:HD21	2.38	0.59
30:D8:26:THR:HB	30:D8:44:VAL:HG22	1.83	0.59
36:1:911:C:H42	39:L2:3:ARG:HD3	1.68	0.59
51:M5:93:LYS:NZ	36:5:2600:C:OP1	156.27	0.59
61:N5:137:ASN:OD1	61:N5:137:ASN:N	2.24	0.59
68:O2:3:SER:HA	68:O2:90:LYS:HB3	3.05	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
70:O4:81:CYS:O	70:O4:83:ASN:N	2.55	0.59
75:O9:4:GLN:HG2	36:5:1588:A:C2	127.25	0.59
79:Q3:84:ARG:NH1	79:Q3:88:GLU:OE2	2.34	0.59
2:S0:30:GLN:HE21	2:S0:33:GLN:HG2	9.87	0.59
36:1:3233:C:H2'	36:1:3234:A:C8	2.37	0.59
36:1:3234:A:H61	36:1:3253:G:H22	1.51	0.59
1:2:780:A:C8	26:D4:8:ARG:HB3	2.37	0.59
1:2:866:G:OP1	15:C3:2:GLY:N	2.34	0.59
56:N0:108:GLN:NE2	36:5:1322:U:O2	292.72	0.59
36:5:138:U:H2'	36:5:139:G:C8	2.38	0.59
23:D1:9:VAL:HG22	23:D1:10:GLU:H	1.92	0.59
52:M6:127:LEU:HD22	56:N0:156:VAL:HG12	1.83	0.59
55:M9:10:LEU:HB3	55:M9:41:ILE:HD12	1.84	0.59
49:M3:50:PRO:HG3	71:O5:118:ILE:HD11	1.84	0.59
71:O5:78:LYS:HG2	71:O5:81:ARG:HD2	3.85	0.59
3:S1:179:SER:HB3	3:S1:183:GLN:HB2	1.84	0.59
1:2:1064:G:O2'	3:S1:204:ILE:O	2.21	0.59
5:S3:137:VAL:HG22	5:S3:151:LYS:HG3	1.84	0.59
36:1:2403:G:N7	36:1:2870:C:H4'	2.18	0.59
1:2:1533:C:H4'	1:2:1539:G:N1	2.17	0.59
36:5:2499:U:H2'	36:5:2500:A:C8	2.37	0.59
1:6:1058:U:H4'	1:6:1059:U:OP1	2.01	0.59
18:C6:42:GLU:HB2	18:C6:45:ARG:HH21	1.67	0.59
21:C9:37:VAL:HG12	21:C9:39:THR:H	4.23	0.59
40:L3:67:PHE:HA	40:L3:70:ARG:HG3	3.96	0.59
41:L4:204:GLY:O	41:L4:246:ARG:NH1	2.35	0.59
44:L7:121:LYS:HE2	44:L7:125:GLU:OE2	2.78	0.59
38:4:155:A:H5'	45:L8:185:ARG:CZ	2.31	0.59
50:M4:121:MET:HE3	36:5:3214:U:H2'	276.70	0.59
54:M8:100:THR:HG23	54:M8:120:GLU:HB3	1.83	0.59
36:1:1320:C:O2	56:N0:115:ARG:NH2	2.35	0.59
59:N3:75:PRO:HG2	59:N3:105:PRO:HD3	1.83	0.59
74:O8:8:ILE:HD12	74:O8:8:ILE:H	2.35	0.59
7:S5:55:ASP:HB3	7:S5:58:LEU:HD12	1.83	0.59
1:2:1347:U:O2	1:2:1516:A:H5''	2.01	0.59
1:2:1594:G:H5'	31:D9:33:LYS:HE3	1.85	0.59
1:2:615:A:H2'	1:2:616:G:H8	1.68	0.59
36:5:2518:C:H2'	36:5:2519:A:C8	2.38	0.59
36:5:3343:G:O2'	36:5:3362:A:N6	2.30	0.59
1:6:482:U:H2'	1:6:483:A:C8	2.38	0.59
13:C1:4:GLU:HG2	13:C1:5:LEU:HG	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:903:U:OP2	16:C4:24:ASN:ND2	2.36	0.59
27:D5:60:VAL:HG13	27:D5:101:TYR:HB2	1.82	0.59
31:D9:5:ASN:HB3	31:D9:7:TRP:NE1	2.18	0.59
44:L7:71:ALA:O	44:L7:73:GLY:N	2.35	0.59
46:L9:117:PHE:O	46:L9:120:ASP:HB2	2.03	0.59
46:L9:89:LYS:HG2	46:L9:145:VAL:HG22	1.85	0.59
70:O4:80:ARG:HG3	70:O4:88:ARG:NH2	3.32	0.59
2:S0:9:LEU:HD23	2:S0:54:TRP:CG	2.37	0.59
4:S2:38:VAL:HG13	4:S2:39:THR:HG23	1.84	0.59
34:SR:112:SER:HB3	34:SR:154:VAL:HG22	1.83	0.59
34:SR:81:LEU:HD11	34:SR:122:ILE:HD13	1.82	0.59
36:1:2768:U:H2'	36:1:2769:A:H8	1.66	0.59
1:2:1087:A:H2'	1:2:1088:A:C8	2.38	0.59
1:2:1381:U:O4	1:2:1382:A:N6	2.36	0.59
1:2:1473:U:H5	7:S5:98:MET:HA	1.68	0.59
1:6:1018:U:H2'	1:6:1019:A:C8	2.38	0.59
15:C3:112:LYS:NZ	1:6:975:C:OP1	277.37	0.59
2:S0:52:LYS:NZ	23:D1:82:VAL:O	2.57	0.59
24:D2:102:VAL:O	24:D2:113:HIS:N	2.36	0.59
39:L2:143:GLU:O	39:L2:145:LYS:N	2.33	0.59
40:L3:212:ASN:O	40:L3:281:LYS:NZ	2.36	0.59
49:M3:120:GLN:HA	49:M3:123:ILE:HG12	1.85	0.59
54:M8:34:THR:HG22	54:M8:49:LEU:HD21	2.23	0.59
59:N3:2:SER:N	59:N3:56:ASP:OD1	6.29	0.59
69:O3:59:VAL:O	69:O3:61:GLY:N	2.35	0.59
2:S0:31:VAL:HG12	2:S0:33:GLN:H	1.67	0.59
9:S7:177:THR:OG1	9:S7:178:GLY:N	2.34	0.59
11:S9:112:GLN:HG3	11:S9:148:VAL:HB	1.84	0.59
36:1:1146:C:H4'	36:1:1331:U:C4	2.38	0.59
36:1:1634:G:N7	63:N7:17:ARG:NH2	2.37	0.59
36:1:1808:G:OP2	63:N7:133:LYS:NZ	2.36	0.59
36:1:2616:C:H3'	36:1:2617:U:O2	2.03	0.59
15:C3:73:ARG:HD3	1:6:859:A:C6	329.72	0.59
16:C4:85:ALA:N	16:C4:119:THR:HG22	2.34	0.59
16:C4:128:LYS:HD2	28:D6:22:ARG:HB3	3.49	0.59
25:D3:44:GLY:H	25:D3:78:LYS:NZ	2.82	0.59
42:L5:136:GLU:CD	42:L5:136:GLU:H	4.80	0.59
42:L5:211:LEU:HB3	42:L5:219:PHE:HB2	2.73	0.59
47:M0:9:TYR:OH	47:M0:99:ILE:HG22	5.06	0.59
62:N6:40:ARG:HG2	62:N6:45:ILE:O	2.02	0.59
62:N6:51:ARG:HG3	62:N6:52:ARG:N	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:SR:108:SER:OG	34:SR:109:ASP:N	2.66	0.59
36:1:829:U:H3	36:1:895:A:H62	1.50	0.59
1:2:1684:U:O2	1:2:1718:G:N2	2.36	0.59
38:4:53:A:H3'	38:4:54:A:H8	1.66	0.59
36:5:1025:A:N7	36:5:1026:A:O2'	2.36	0.59
64:N8:21:ARG:HD2	36:5:1369:A:H5''	184.78	0.59
36:5:91:G:OP2	36:5:93:C:N4	2.30	0.59
1:6:230:C:H42	1:6:235:G:H1	1.51	0.59
16:C4:104:ALA:HA	16:C4:107:ARG:HB3	3.17	0.59
17:C5:118:GLU:O	20:C8:122:HIS:N	2.93	0.59
20:C8:30:TYR:HE2	20:C8:40:ARG:HH11	1.64	0.59
40:L3:92:TYR:HB2	40:L3:157:VAL:HG13	3.73	0.59
36:1:1381:A:OP1	41:L4:197:ARG:NH1	2.36	0.59
55:M9:21:LYS:HE3	55:M9:55:VAL:HA	4.45	0.59
63:N7:95:VAL:HG21	63:N7:113:VAL:HG11	2.25	0.59
64:N8:74:ASN:HB3	64:N8:115:LYS:H	1.68	0.59
36:1:975:C:H2'	36:1:976:U:H6	1.68	0.59
1:2:918:U:H2'	1:2:919:A:C8	2.38	0.59
36:5:1635:G:N2	36:5:1638:A:OP2	2.33	0.59
36:5:2996:U:OP1	36:5:2996:U:H4'	2.02	0.59
36:5:3378:C:H2'	36:5:3379:C:H6	1.68	0.59
48:M1:137:ARG:HG2	37:7:28:C:H5''	308.95	0.59
19:C7:27:ASP:O	19:C7:31:ASN:ND2	2.36	0.59
20:C8:82:PRO:O	20:C8:84:TRP:N	2.32	0.59
23:D1:74:GLN:HE22	23:D1:83:TRP:H	1.50	0.59
30:D8:36:THR:OG1	30:D8:37:SER:N	2.36	0.59
41:L4:44:LYS:O	41:L4:47:ARG:HD2	2.63	0.59
41:L4:89:ALA:O	41:L4:91:GLY:N	2.33	0.59
44:L7:160:ARG:HG3	44:L7:203:TRP:CD2	2.37	0.59
56:N0:74:ASN:HD21	56:N0:95:ARG:HH11	1.50	0.59
36:1:654:C:OP1	68:O2:27:ARG:NH2	2.36	0.59
6:S4:160:VAL:HG13	6:S4:169:ILE:HG23	2.03	0.59
6:S4:187:ARG:NH2	1:6:753:A:N7	374.42	0.59
35:SM:48:ARG:HH12	36:5:1017:C:H5''	338.33	0.59
34:SR:58:VAL:HG23	34:SR:59:ARG:HG3	1.84	0.59
36:1:1080:A:OP2	42:L5:140:ARG:NH2	2.36	0.58
36:1:1721:U:OP2	55:M9:124:TYR:OH	2.13	0.58
36:1:2112:U:H4'	36:1:2113:A:H5'	1.84	0.58
36:1:31:C:OP2	51:M5:188:ARG:NH2	2.36	0.58
37:3:11:A:H4'	37:3:13:A:C8	2.38	0.58
55:M9:9:ARG:HH21	36:5:1603:A:H5'	109.88	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:17:ARG:NH2	36:5:1634:G:N7	197.41	0.58
36:5:2396:G:OP1	36:5:2397:A:H4'	2.03	0.58
57:N1:68:THR:OG1	36:5:2737:C:H4'	224.40	0.58
21:C9:128:GLY:HA2	21:C9:131:ASP:HB2	3.49	0.58
27:D5:88:ILE:HA	27:D5:104:ALA:HB2	1.85	0.58
39:L2:59:ALA:HB2	39:L2:78:ALA:HB2	1.84	0.58
41:L4:255:PHE:HA	41:L4:258:LEU:HD22	2.39	0.58
51:M5:122:ASN:OD1	51:M5:123:GLN:N	2.36	0.58
55:M9:102:LEU:HD22	55:M9:138:LEU:HD12	1.85	0.58
67:O1:15:ASN:O	67:O1:19:ARG:NH1	3.56	0.58
36:1:362:U:O4	73:O7:24:ARG:NH2	2.36	0.58
6:S4:52:LEU:HB3	6:S4:54:TYR:HD2	1.68	0.58
34:SR:133:VAL:O	34:SR:141:LEU:N	2.34	0.58
34:SR:198:ASN:O	34:SR:215:GLY:HA3	2.17	0.58
36:1:1350:A:H2'	36:1:1351:U:H3'	1.85	0.58
36:5:1783:U:H2'	36:5:1784:G:C8	2.38	0.58
47:M0:157:TYR:CD1	36:5:2836:C:H4'	311.71	0.58
36:5:528:U:H2'	36:5:529:A:C8	2.37	0.58
75:O9:21:ARG:NH2	38:8:51:G:OP2	76.66	0.58
5:S3:23:GLU:CD	12:C0:61:TRP:HE1	2.06	0.58
16:C4:84:ARG:HB3	16:C4:118:VAL:HG23	3.22	0.58
41:L4:302:ALA:HB2	54:M8:39:ARG:NH2	2.69	0.58
42:L5:77:ALA:O	42:L5:108:ARG:NH1	2.36	0.58
45:L8:121:SER:O	45:L8:123:GLN:N	2.36	0.58
74:O8:32:ASN:O	74:O8:34:ALA:N	2.36	0.58
3:S1:99:ASN:OD1	3:S1:100:PHE:N	2.36	0.58
4:S2:73:LEU:HG	4:S2:76:LEU:HD13	1.84	0.58
1:2:1145:U:O2'	4:S2:89:GLN:O	2.14	0.58
36:1:1315:U:OP2	52:M6:44:SER:OG	2.17	0.58
1:2:1160:A:H2'	1:2:1161:C:C6	2.38	0.58
1:2:1208:A:N1	1:2:1455:G:N2	2.48	0.58
36:5:1239:C:H42	36:5:1249:G:H1	1.51	0.58
36:5:2927:C:H2'	36:5:2928:C:C6	2.38	0.58
61:N5:56:ARG:NH2	38:8:135:G:OP2	81.97	0.58
38:8:149:A:H2'	38:8:150:G:C8	2.38	0.58
40:L3:4:ARG:HD3	40:L3:7:GLU:HA	1.85	0.58
41:L4:93:MET:H	41:L4:93:MET:HE2	1.67	0.58
42:L5:180:PHE:HB3	42:L5:195:LEU:HD13	1.94	0.58
46:L9:173:ARG:NH2	36:5:2898:G:OP2	330.08	0.58
50:M4:48:GLY:HA3	50:M4:53:VAL:HG13	2.14	0.58
63:N7:115:LYS:NZ	63:N7:119:GLU:OE2	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S1:67:GLU:OE1	16:C4:82:LYS:NZ	2.37	0.58
8:S6:13:GLN:OE1	1:6:151:G:N2	310.87	0.58
9:S7:162:ILE:HG22	9:S7:165:LYS:HD2	1.85	0.58
34:SR:149:ASP:HB2	34:SR:175:ASP:HB3	1.89	0.58
36:1:1596:C:H2'	36:1:1597:C:C6	2.39	0.58
36:1:1722:U:OP1	55:M9:100:ARG:HD3	2.03	0.58
36:1:519:A:OP1	56:N0:62:ASN:ND2	2.33	0.58
36:1:692:A:OP1	51:M5:201:ARG:NH2	2.35	0.58
42:L5:23:ARG:NH2	36:5:2703:A:OP2	283.58	0.58
36:5:3299:A:N6	36:5:3315:G:H1	1.99	0.58
42:L5:41:LYS:HD2	57:N1:93:VAL:HG11	1.85	0.58
43:L6:129:GLU:N	43:L6:129:GLU:OE1	2.37	0.58
47:M0:61:SER:OG	47:M0:63:GLU:HG2	4.05	0.58
59:N3:129:VAL:O	59:N3:133:SER:OG	3.43	0.58
78:Q2:63:LYS:HD2	78:Q2:87:ARG:NH1	2.19	0.58
7:S5:41:LYS:NZ	7:S5:67:PRO:O	3.76	0.58
35:SM:50:ASN:N	35:SM:50:ASN:OD1	4.07	0.58
1:2:473:A:H5'	1:2:769:A:H1'	1.85	0.58
36:5:2709:C:H2'	36:5:2710:C:C6	2.38	0.58
36:5:59:G:H2'	38:8:33:A:O2'	2.04	0.58
51:M5:172:ARG:NH2	36:5:63:A:OP1	103.72	0.58
38:8:83:C:H4'	38:8:85:G:N3	2.18	0.58
13:C1:22:ASN:HD22	13:C1:25:VAL:HG23	1.68	0.58
33:E1:108:VAL:HG12	33:E1:114:VAL:HG13	3.92	0.58
33:E1:90:LYS:HB2	33:E1:93:HIS:CE1	12.71	0.58
8:S6:94:ARG:HH21	1:6:407:A:H5'	289.50	0.58
9:S7:143:LEU:HD23	9:S7:149:ILE:HD13	4.64	0.58
36:1:1720:U:OP2	55:M9:120:TYR:OH	2.14	0.58
35:SM:34:LYS:NZ	36:1:2707:C:OP1	2.37	0.58
36:1:386:A:C5	36:1:387:A:H1'	2.39	0.58
1:2:1257:U:H2'	12:C0:2:LEU:HD12	1.85	0.58
36:5:22:G:H1'	38:8:104:A:N3	2.18	0.58
4:S2:143:TYR:O	24:D2:98:GLN:NE2	3.35	0.58
25:D3:126:LYS:HG3	25:D3:131:SER:HA	1.85	0.58
46:L9:163:GLN:OE1	46:L9:166:ARG:NH1	3.60	0.58
58:N2:50:LEU:HD22	58:N2:54:VAL:HB	4.75	0.58
67:O1:31:ARG:O	67:O1:35:GLU:HB2	2.38	0.58
69:O3:26:ASN:HA	69:O3:88:ASN:OD1	2.04	0.58
6:S4:191:ARG:HD3	6:S4:245:LYS:HB2	1.85	0.58
7:S5:79:ASN:OD1	7:S5:83:ARG:NH2	3.84	0.58
10:S8:122:GLY:N	10:S8:157:GLU:OE2	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:SM:49:LYS:N	36:1:1019:G:OP1	2.37	0.58
36:1:1761:C:O2'	36:1:1762:C:OP2	2.22	0.58
36:5:2861:U:H2'	36:5:2862:U:O4'	2.03	0.58
13:C1:98:ASN:ND2	13:C1:98:ASN:O	2.37	0.58
22:D0:24:ILE:HA	22:D0:116:VAL:HG12	4.87	0.58
22:D0:71:PRO:O	22:D0:72:ASN:ND2	5.18	0.58
33:E1:90:LYS:HB2	33:E1:93:HIS:HE1	13.12	0.58
40:L3:168:LYS:O	40:L3:319:ASN:ND2	2.84	0.58
50:M4:17:VAL:HG21	50:M4:74:ARG:HB2	1.85	0.58
51:M5:73:ARG:HG2	51:M5:75:VAL:HG13	2.06	0.58
53:M7:177:ALA:HA	53:M7:180:LYS:HD2	1.84	0.58
54:M8:65:SER:OG	54:M8:90:ASP:OD2	2.44	0.58
61:N5:135:ILE:HA	61:N5:138:ARG:HB3	2.78	0.58
64:N8:24:LYS:HD2	64:N8:26:ARG:HE	1.68	0.58
68:O2:71:HIS:CE1	68:O2:118:LYS:HD3	2.38	0.58
79:Q3:59:CYS:O	79:Q3:61:LYS:N	2.36	0.58
1:2:14:C:H5''	4:S2:203:LYS:HD2	1.85	0.58
4:S2:222:TYR:HE2	23:D1:11:LEU:HD12	3.49	0.58
5:S3:142:LEU:HD13	5:S3:182:LEU:HD21	1.85	0.58
6:S4:246:LEU:H	6:S4:246:LEU:HD12	1.67	0.58
6:S4:251:GLU:HB3	6:S4:255:ARG:NH1	2.19	0.58
9:S7:50:ASP:OD1	9:S7:50:ASP:N	2.85	0.58
10:S8:43:ILE:H	1:6:260:U:H5	273.65	0.58
36:1:398:A:O2'	36:1:1416:C:OP1	2.18	0.58
36:1:1688:U:H2'	36:1:1689:U:C6	2.39	0.58
36:1:2748:A:O3'	42:L5:48:LYS:NZ	2.37	0.58
1:2:1266:U:H2'	1:2:1267:G:C8	2.39	0.58
1:2:323:A:OP2	10:S8:10:LYS:HG3	2.03	0.58
38:4:151:C:O2'	38:4:153:U:OP2	2.21	0.58
36:5:1561:G:O6	36:5:1578:C:N4	2.37	0.58
6:S4:187:ARG:HH22	1:6:753:A:H62	375.51	0.58
7:S5:25:LEU:HB3	18:C6:27:GLY:HA3	3.03	0.58
30:D8:12:VAL:HA	30:D8:30:VAL:HG12	1.85	0.58
31:D9:24:CYS:HB3	31:D9:42:CYS:SG	4.02	0.58
39:L2:238:ILE:H	39:L2:238:ILE:HD12	1.67	0.58
41:L4:23:PRO:O	41:L4:25:VAL:N	2.37	0.58
42:L5:120:LYS:O	42:L5:248:ARG:NH2	2.42	0.58
44:L7:88:ARG:HD2	44:L7:90:LYS:O	2.20	0.58
45:L8:84:ARG:HH22	45:L8:181:LYS:HZ3	1.51	0.58
49:M3:170:LEU:HD22	72:O6:9:ILE:HD11	1.85	0.58
52:M6:27:LEU:HD22	52:M6:101:ARG:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:N0:60:SER:OG	56:N0:62:ASN:OD1	2.21	0.58
73:O7:29:VAL:O	73:O7:32:LYS:NZ	4.13	0.58
73:O7:45:ARG:NH2	36:5:361:A:O3'	124.11	0.58
2:S0:139:VAL:HG13	2:S0:141:ILE:HG13	1.86	0.58
4:S2:88:LYS:HB3	4:S2:95:ARG:HD2	4.69	0.58
36:1:1051:U:H4'	57:N1:19:PHE:CD2	2.38	0.58
36:1:2996:U:OP1	36:1:2996:U:H4'	2.02	0.58
1:2:381:C:O2'	1:2:755:A:N1	2.36	0.58
36:5:1355:A:H4'	36:5:1356:U:O5'	2.04	0.58
45:L8:157:VAL:HG13	36:5:147:U:C4	127.36	0.58
36:5:379:C:H2'	36:5:380:U:H6	1.69	0.58
13:C1:83:THR:HA	13:C1:111:VAL:H	2.37	0.58
15:C3:92:ILE:O	15:C3:96:VAL:HG23	2.03	0.58
1:2:375:U:OP1	25:D3:23:ARG:NH2	2.37	0.58
25:D3:46:SER:OG	25:D3:48:HIS:O	2.21	0.58
36:1:2415:C:OP1	39:L2:2:GLY:N	2.36	0.58
42:L5:257:GLU:N	42:L5:257:GLU:OE2	4.26	0.58
49:M3:108:ILE:HG22	49:M3:112:ASN:HD21	1.68	0.58
51:M5:73:ARG:O	51:M5:75:VAL:N	3.79	0.58
72:O6:76:ARG:O	36:5:293:C:O2'	152.12	0.58
2:S0:124:THR:HA	2:S0:146:LEU:HB2	1.84	0.58
35:SM:46:LYS:HA	36:5:1018:G:H4'	324.28	0.58
1:2:1445:G:H1	33:E1:89:LYS:HA	1.69	0.58
1:2:1535:U:O2'	1:2:1536:G:N3	2.36	0.58
1:6:213:A:H61	1:6:252:U:H3	1.52	0.58
1:6:340:U:H2'	1:6:341:A:C8	2.39	0.58
15:C3:101:HIS:HA	15:C3:104:ARG:HE	1.69	0.58
20:C8:14:ILE:HD12	20:C8:23:ASP:HA	5.96	0.58
24:D2:83:ILE:HD12	24:D2:122:SER:HB2	4.71	0.58
40:L3:167:ARG:HH11	40:L3:167:ARG:HB2	1.68	0.58
41:L4:74:ILE:HD13	41:L4:88:GLY:HA2	1.86	0.58
42:L5:196:ARG:HA	42:L5:199:ILE:HD12	2.40	0.58
49:M3:22:VAL:HG13	51:M5:197:LEU:HB3	1.85	0.58
55:M9:43:LYS:O	55:M9:47:ASN:N	3.29	0.58
36:1:1233:G:H22	36:1:1255:C:H42	1.52	0.57
36:1:3086:A:OP1	40:L3:367:LYS:NZ	2.22	0.57
1:2:1081:A:O2'	1:2:1082:C:O5'	2.15	0.57
1:2:733:A:H4'	1:2:734:A:C5	2.39	0.57
37:3:65:G:OP1	47:M0:204:GLY:N	2.27	0.57
36:5:260:C:H2'	36:5:261:U:C6	2.38	0.57
27:D5:77:ARG:NH2	1:6:1534:G:N7	348.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:560:U:H2'	1:6:561:G:C8	2.39	0.57
17:C5:30:THR:HG23	17:C5:86:VAL:HG11	2.95	0.57
19:C7:75:GLU:O	19:C7:79:GLU:HG2	2.04	0.57
39:L2:221:LYS:O	36:5:2245:C:H4'	218.62	0.57
45:L8:45:ASN:ND2	45:L8:47:SER:HB3	2.18	0.57
46:L9:37:ASN:O	46:L9:39:LYS:N	2.36	0.57
50:M4:38:ILE:HD11	56:N0:150:PHE:CE2	2.39	0.57
67:O1:44:MET:HB3	67:O1:77:ARG:NH1	4.06	0.57
36:1:623:U:O3'	69:O3:86:ARG:NH2	2.36	0.57
4:S2:168:ARG:NE	1:6:1098:U:OP2	383.48	0.57
10:S8:89:GLU:O	10:S8:93:THR:OG1	2.11	0.57
36:1:2355:G:H4'	53:M7:139:TYR:CE2	2.39	0.57
36:1:2623:G:H2'	36:1:2624:G:C8	2.38	0.57
1:2:1031:U:H4'	1:2:1032:G:OP2	2.04	0.57
36:5:2899:C:O2'	36:5:2901:G:OP2	2.17	0.57
17:C5:68:PRO:HG2	17:C5:71:GLU:HG3	1.85	0.57
28:D6:15:ARG:O	28:D6:17:HIS:N	2.37	0.57
36:1:63:A:H5''	51:M5:174:ILE:HG21	1.86	0.57
51:M5:190:THR:O	51:M5:194:GLN:HG2	2.29	0.57
67:O1:41:LYS:NZ	67:O1:47:ASP:OD1	2.36	0.57
4:S2:140:ARG:HG2	4:S2:155:ALA:HB2	3.46	0.57
7:S5:203:LYS:O	7:S5:205:SER:N	3.54	0.57
34:SR:164:ASP:O	34:SR:166:SER:N	3.17	0.57
36:1:1805:C:H2'	36:1:1806:A:H8	1.68	0.57
36:1:2620:G:H1	84:C:74:CH:H41	1.52	0.57
36:1:2812:C:H2'	36:1:2813:A:C8	2.40	0.57
1:2:1222:C:H42	1:2:1261:G:H1	1.51	0.57
1:2:1335:U:H3	1:2:1416:G:H1	1.50	0.57
36:5:2404:A:N7	36:5:2872:A:N6	2.52	0.57
36:5:327:A:H2'	36:5:328:U:C6	2.39	0.57
36:5:561:C:H2'	36:5:562:C:C6	2.38	0.57
19:C7:52:GLY:HA3	1:6:1389:C:O2'	422.38	0.57
19:C7:108:ASP:HA	19:C7:111:LYS:HE3	4.80	0.57
6:S4:59:ARG:HH12	26:D4:87:PRO:HG3	1.68	0.57
39:L2:27:ALA:O	39:L2:128:ARG:NH2	2.38	0.57
40:L3:19:ARG:HB3	40:L3:232:ARG:NH1	2.42	0.57
40:L3:284:ARG:NH2	40:L3:295:ALA:O	2.65	0.57
46:L9:163:GLN:O	46:L9:166:ARG:HG3	3.62	0.57
47:M0:93:PRO:HB2	47:M0:125:LEU:HB2	3.59	0.57
64:N8:125:VAL:HG21	64:N8:138:ILE:HD13	1.87	0.57
71:O5:89:ARG:HD2	38:8:38:U:C4	68.23	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
78:Q2:77:CYS:SG	78:Q2:79:THR:OG1	2.86	0.57
3:S1:36:SER:HB3	3:S1:41:ARG:HH21	7.73	0.57
5:S3:59:LEU:HD23	5:S3:66:ILE:HG21	7.87	0.57
72:O6:14:GLY:HA2	36:5:73:C:OP1	107.72	0.57
1:6:1041:G:H2'	1:6:1042:G:C8	2.39	0.57
1:6:104:A:N6	1:6:308:C:H5'	2.10	0.57
1:6:751:G:H2'	1:6:752:A:C8	2.39	0.57
1:2:1565:C:OP1	20:C8:41:ARG:HG3	2.04	0.57
25:D3:102:VAL:HG12	25:D3:127:VAL:HA	2.28	0.57
30:D8:32:PHE:HE2	30:D8:38:ARG:HB3	1.69	0.57
39:L2:114:SER:HB2	39:L2:169:ILE:HG12	2.93	0.57
42:L5:211:LEU:O	42:L5:215:ASP:N	3.82	0.57
44:L7:27:ALA:HA	44:L7:30:ARG:HB3	1.86	0.57
46:L9:90:MET:HG2	46:L9:181:VAL:HG22	1.86	0.57
47:M0:29:SER:OG	47:M0:30:LYS:N	2.36	0.57
62:N6:35:LEU:HD21	62:N6:48:LEU:HD12	1.86	0.57
6:S4:2:ALA:O	6:S4:3:ARG:HB2	2.34	0.57
6:S4:95:THR:OG1	6:S4:95:THR:O	3.46	0.57
8:S6:57:ASP:O	8:S6:59:GLN:N	3.65	0.57
34:SR:255:ALA:HA	34:SR:260:ILE:HG12	1.85	0.57
1:2:1059:U:O2'	1:2:1060:U:N3	2.37	0.57
67:O1:28:ARG:NH1	36:5:3056:U:O2	184.36	0.57
43:L6:78:ARG:NH1	36:5:3272:C:OP2	247.01	0.57
1:6:485:A:C6	1:6:486:G:H1'	2.39	0.57
1:6:755:A:H2'	1:6:756:A:C8	2.38	0.57
30:D8:38:ARG:HH11	30:D8:40:ILE:HD11	1.69	0.57
48:M1:155:THR:O	48:M1:159:THR:HG23	5.30	0.57
36:1:1874:A:N7	55:M9:20:ARG:NH1	2.53	0.57
65:N9:14:ARG:HH12	65:N9:18:ARG:HD2	1.70	0.57
67:O1:25:PHE:HB3	67:O1:65:LYS:HG2	1.86	0.57
73:O7:75:LYS:HD3	36:5:181:U:H4'	50.89	0.57
3:S1:26:ARG:NH1	3:S1:49:ASN:OD1	2.38	0.57
6:S4:146:THR:HG21	1:6:123:G:H21	340.61	0.57
36:1:2707:C:H2'	36:1:2708:C:C6	2.40	0.57
36:1:3001:C:OP1	40:L3:120:LYS:NZ	2.31	0.57
36:1:422:A:C2	36:1:2363:A:H4'	2.40	0.57
36:5:1244:A:H5''	36:5:1271:A:H4'	1.86	0.57
36:5:1493:G:OP2	36:5:1493:G:N2	2.36	0.57
36:5:70:A:N1	36:5:313:A:O2'	2.36	0.57
25:D3:97:ASP:O	25:D3:100:ASP:HB2	3.67	0.57
30:D8:22:ARG:NH1	1:6:1619:C:O2	339.80	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:M7:122:ALA:HB3	53:M7:143:PRO:HB2	1.91	0.57
62:N6:5:SER:C	62:N6:7:ASP:H	2.49	0.57
2:S0:17:LEU:HD23	2:S0:172:LEU:HD13	1.86	0.57
6:S4:73:ASP:OD2	6:S4:122:LYS:NZ	2.24	0.57
34:SR:33:LEU:HB2	34:SR:47:LEU:HD11	1.85	0.57
36:1:1095:U:O2	57:N1:128:LEU:N	2.38	0.57
36:1:1942:U:OP2	55:M9:74:ARG:NH1	2.38	0.57
1:2:26:A:O2'	1:2:27:U:O5'	2.23	0.57
36:5:1110:U:H2'	36:5:1111:U:C6	2.40	0.57
36:5:2209:U:H4'	36:5:2210:G:OP1	2.04	0.57
1:6:1209:C:H42	1:6:1454:G:H1	1.52	0.57
1:6:427:C:O2'	1:6:459:G:N3	2.28	0.57
1:6:691:C:P	1:6:696:C:H41	2.27	0.57
12:C0:15:LEU:HD13	12:C0:21:VAL:HG23	1.87	0.57
27:D5:38:HIS:ND1	27:D5:71:ILE:O	7.19	0.57
40:L3:308:MET:HB2	40:L3:363:SER:HB2	1.87	0.57
41:L4:131:VAL:HG12	41:L4:134:LEU:H	2.53	0.57
44:L7:86:VAL:HG21	44:L7:127:LEU:HD21	2.29	0.57
45:L8:94:PHE:HB3	45:L8:189:LEU:HD13	1.87	0.57
71:O5:84:LYS:O	73:O7:73:ARG:NH2	3.73	0.57
74:O8:29:LYS:NZ	74:O8:29:LYS:O	2.25	0.57
3:S1:88:VAL:HG11	3:S1:96:LEU:HD12	1.86	0.57
36:1:1877:U:H5''	36:1:1878:G:O4'	2.03	0.57
36:1:289:A:O2'	51:M5:93:LYS:O	2.23	0.57
36:1:3316:A:H2	36:1:3389:U:H5'	1.70	0.57
57:N1:43:LYS:HD2	36:5:992:A:H5''	256.31	0.57
9:S7:115:SER:O	1:6:856:A:N6	359.38	0.57
22:D0:22:ILE:HD12	22:D0:118:VAL:HA	1.87	0.57
43:L6:81:ALA:O	43:L6:84:VAL:HG12	2.04	0.57
44:L7:158:LYS:HG2	44:L7:203:TRP:HH2	1.69	0.57
40:L3:57:VAL:HG21	60:N4:15:PRO:HG2	1.87	0.57
3:S1:175:GLU:HG3	3:S1:193:ILE:HG23	1.86	0.57
3:S1:36:SER:HA	3:S1:41:ARG:HE	4.08	0.57
6:S4:72:VAL:HB	6:S4:77:ARG:HG3	4.18	0.57
11:S9:119:ALA:HA	11:S9:124:HIS:HD2	2.84	0.57
36:1:1185:C:OP1	50:M4:42:LYS:HD3	2.05	0.57
36:1:201:A:H2'	36:1:202:G:H8	1.69	0.57
36:1:3275:U:O4'	69:O3:66:VAL:HG21	2.05	0.57
1:2:800:U:H2'	1:2:801:G:H8	1.70	0.57
1:2:926:A:C2	16:C4:125:SER:HB3	2.39	0.57
36:5:2209:U:H1'	36:5:2210:G:H5''	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:C3:105:ASN:HB3	1:6:879:G:O2'	273.78	0.57
15:C3:67:THR:O	15:C3:67:THR:OG1	2.63	0.57
15:C3:91:LEU:HD12	15:C3:125:LEU:HD12	1.86	0.57
17:C5:22:LEU:HA	17:C5:25:LEU:HD12	2.69	0.57
44:L7:229:PHE:C	44:L7:229:PHE:HD1	2.44	0.57
49:M3:124:ILE:HD12	49:M3:126:PHE:HE1	2.06	0.57
42:L5:34:LYS:HD2	57:N1:30:TYR:CE2	2.40	0.57
64:N8:73:LEU:HD23	64:N8:112:ILE:HD12	1.87	0.57
2:S0:146:LEU:HD23	2:S0:160:ILE:HB	5.15	0.57
3:S1:145:LYS:HG3	3:S1:149:GLN:HB3	5.57	0.57
4:S2:41:LEU:HD11	4:S2:56:ILE:HG21	1.87	0.57
5:S3:192:PRO:HB2	5:S3:201:ALA:HA	2.52	0.57
36:1:645:A:N6	36:1:2869:U:OP1	2.32	0.57
36:1:3094:A:H2'	36:1:3095:U:H6	1.70	0.57
36:1:417:A:H2'	36:1:418:A:C8	2.40	0.57
49:M3:39:ARG:NH1	36:5:107:A:OP1	74.05	0.57
76:Q0:114:LYS:NZ	36:5:3107:U:OP1	300.04	0.57
1:6:1151:A:O2'	1:6:1766:A:N7	2.28	0.57
18:C6:58:ASP:O	18:C6:60:PHE:N	2.38	0.57
19:C7:82:ASP:O	19:C7:83:GLN:NE2	2.37	0.57
21:C9:30:VAL:O	21:C9:32:GLY:N	2.38	0.57
24:D2:10:ALA:HB1	24:D2:27:ILE:HD12	1.87	0.57
29:D7:72:LYS:NZ	1:6:1063:U:OP1	340.10	0.57
33:E1:134:ASN:OD1	33:E1:134:ASN:N	4.27	0.57
41:L4:329:PRO:HB2	44:L7:45:LEU:HD12	2.34	0.57
49:M3:35:ARG:NH1	36:5:685:G:OP2	83.63	0.57
62:N6:106:ILE:HG21	62:N6:109:LEU:HD23	2.38	0.57
68:O2:11:LYS:HB3	68:O2:14:THR:HG22	1.87	0.57
3:S1:137:ILE:HD12	3:S1:172:LEU:HD22	1.87	0.57
3:S1:176:VAL:HG13	3:S1:184:LEU:HD21	1.87	0.57
1:2:816:G:N2	9:S7:110:GLN:OE1	2.31	0.57
1:2:4:C:O2'	11:S9:17:ARG:NH1	2.38	0.57
1:2:1085:G:N2	1:2:1088:A:OP2	2.38	0.56
36:5:182:U:H2'	36:5:183:G:C8	2.40	0.56
36:5:2413:A:H2'	36:5:2414:G:H8	1.69	0.56
36:5:2712:U:H2'	36:5:2713:U:C6	2.39	0.56
36:5:3317:U:H4'	36:5:3318:G:O5'	2.05	0.56
19:C7:48:ASN:ND2	1:6:1388:A:H5''	430.62	0.56
1:6:1413:U:H4'	1:6:1414:U:OP2	2.05	0.56
27:D5:77:ARG:NH1	1:6:1533:C:OP2	352.91	0.56
38:8:44:A:H2'	38:8:45:C:C6	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:C9:28:LEU:HD23	21:C9:111:ILE:HD11	7.69	0.56
16:C4:112:ILE:H	28:D6:57:SER:HA	2.04	0.56
46:L9:117:PHE:CE1	46:L9:165:CYS:HB3	2.71	0.56
46:L9:49:ASN:O	46:L9:49:ASN:ND2	2.37	0.56
53:M7:29:THR:O	53:M7:119:VAL:HG11	2.04	0.56
36:1:3174:A:OP1	69:O3:97:SER:OG	2.22	0.56
3:S1:143:THR:HA	3:S1:207:LEU:HA	1.88	0.56
3:S1:164:ILE:HD13	3:S1:207:LEU:HD11	4.22	0.56
8:S6:152:ASP:OD1	8:S6:154:ARG:NH1	5.54	0.56
36:1:732:C:H42	36:1:737:G:H1	1.53	0.56
36:5:2112:U:H4'	36:5:2113:A:O5'	2.05	0.56
1:6:1280:C:H2'	1:6:1281:G:H8	1.70	0.56
6:S4:187:ARG:NH2	1:6:753:A:H62	374.94	0.56
6:S4:106:LYS:NZ	1:6:788:A:OP1	396.94	0.56
1:6:973:A:H2'	1:6:974:A:H8	1.69	0.56
14:C2:55:GLY:HA2	14:C2:85:LYS:HE3	1.85	0.56
15:C3:99:ARG:O	15:C3:103:GLU:HG2	2.05	0.56
28:D6:64:LEU:HD23	28:D6:65:PRO:HD2	4.93	0.56
42:L5:35:ARG:HG2	36:5:2749:G:O2'	249.23	0.56
52:M6:18:ARG:O	52:M6:22:VAL:HG12	2.62	0.56
38:4:131:A:H5''	61:N5:93:TYR:CE2	2.40	0.56
2:S0:55:GLU:OE2	23:D1:80:LYS:N	3.12	0.56
6:S4:151:ASP:HB3	6:S4:154:ILE:HG13	1.87	0.56
34:SR:219:GLU:HA	34:SR:235:SER:HA	2.88	0.56
36:1:763:G:H2'	36:1:764:U:O4'	2.05	0.56
36:1:838:G:O6	79:Q3:4:ARG:NH2	2.38	0.56
1:2:1290:U:H2'	1:2:1291:G:C8	2.40	0.56
36:5:2518:C:H2'	36:5:2519:A:H8	1.70	0.56
36:5:394:G:N2	36:5:396:A:H3'	2.20	0.56
1:6:819:G:O2'	1:6:821:U:OP2	2.23	0.56
37:7:2:G:O2'	37:7:23:A:N1	2.37	0.56
62:N6:116:LYS:NZ	38:8:84:C:N3	29.88	0.56
23:D1:41:GLU:O	23:D1:44:ARG:NH2	2.38	0.56
43:L6:89:THR:HG21	50:M4:115:PHE:HB2	1.88	0.56
53:M7:178:ALA:HA	53:M7:181:ARG:HB3	1.87	0.56
67:O1:70:ARG:HE	67:O1:102:LYS:NZ	6.59	0.56
49:M3:177:LYS:HA	72:O6:11:LEU:HD22	2.71	0.56
72:O6:26:ILE:HD13	36:5:155:G:H1'	87.54	0.56
2:S0:64:ILE:HG23	2:S0:73:VAL:HG11	1.87	0.56
9:S7:8:ILE:HG23	9:S7:42:GLN:HA	1.85	0.56
9:S7:58:LEU:HB2	9:S7:90:VAL:HG22	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1642:A:N3	36:1:1822:C:O2'	2.37	0.56
36:1:3350:C:O2'	36:1:3351:U:O5'	2.23	0.56
1:6:26:A:OP2	1:6:26:A:H3'	2.04	0.56
1:6:60:U:H5'	1:6:455:C:H42	1.70	0.56
37:7:91:G:H2'	37:7:92:A:H8	1.70	0.56
13:C1:80:MET:HB2	13:C1:83:THR:HG23	1.88	0.56
24:D2:8:ALA:HA	24:D2:74:VAL:HG11	1.85	0.56
36:1:805:G:H1'	41:L4:73:ARG:HH11	1.69	0.56
44:L7:140:SER:O	44:L7:144:ILE:HG13	2.74	0.56
43:L6:175:LYS:HD3	50:M4:111:ALA:HA	1.85	0.56
50:M4:23:ILE:HG21	50:M4:28:SER:HB2	1.87	0.56
61:N5:38:LEU:HD11	61:N5:40:LEU:HD13	1.86	0.56
69:O3:45:LEU:HD13	69:O3:71:VAL:HG12	4.97	0.56
36:1:3275:U:H5'	69:O3:68:TRP:CZ2	2.37	0.56
36:1:1491:A:N7	75:O9:2:ALA:HB3	2.20	0.56
4:S2:140:ARG:HB3	4:S2:221:THR:HB	2.60	0.56
36:1:1330:A:OP2	69:O3:19:SER:HB3	2.05	0.56
36:1:688:G:H2'	36:1:690:A:C8	2.40	0.56
36:1:914:A:C5	39:L2:199:THR:HG21	2.40	0.56
1:2:1067:C:H2'	1:2:1068:C:C6	2.41	0.56
1:2:1796:C:O5'	28:D6:5:ARG:NH1	2.39	0.56
1:2:170:U:H6	1:2:267:U:HO2'	1.54	0.56
36:5:1596:C:O2'	36:5:1696:A:N3	2.35	0.56
36:5:1764:U:H3'	36:5:1765:U:H5''	1.86	0.56
55:M9:74:ARG:NH1	36:5:1942:U:OP2	209.08	0.56
36:5:2101:C:H2'	36:5:2102:U:C6	2.41	0.56
67:O1:43:HIS:HE1	36:5:3323:A:H1'	169.14	0.56
1:6:906:A:H2	1:6:998:A:HO2'	1.50	0.56
15:C3:87:ASP:HB3	15:C3:125:LEU:HD11	4.15	0.56
26:D4:7:ILE:HG23	26:D4:27:VAL:HG22	1.87	0.56
42:L5:52:VAL:HA	42:L5:147:ASP:HB3	1.95	0.56
43:L6:18:LEU:H	43:L6:18:LEU:HD22	1.70	0.56
36:1:1334:U:O2'	44:L7:151:ARG:NH2	2.39	0.56
45:L8:78:PHE:C	45:L8:80:TYR:H	2.09	0.56
56:N0:137:ARG:HG2	56:N0:139:TYR:CZ	2.41	0.56
57:N1:130:ARG:O	36:5:1098:A:O2'	256.22	0.56
61:N5:105:VAL:HG11	61:N5:126:LEU:HD13	2.01	0.56
68:O2:97:ALA:HB3	68:O2:100:ILE:HG12	1.86	0.56
69:O3:45:LEU:HA	69:O3:71:VAL:HG12	1.86	0.56
7:S5:90:ILE:O	7:S5:94:THR:HG23	2.05	0.56
8:S6:13:GLN:CD	1:6:151:G:H21	310.50	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:S6:135:PRO:HB2	8:S6:141:ILE:HG12	1.87	0.56
34:SR:50:ASP:O	34:SR:52:GLN:N	2.38	0.56
34:SR:80:ALA:O	34:SR:92:TRP:N	2.85	0.56
36:5:209:A:H1'	36:5:212:G:H22	1.69	0.56
36:5:2413:A:H2'	36:5:2414:G:C8	2.40	0.56
1:6:1229:G:O2'	1:6:1255:G:N2	2.39	0.56
16:C4:35:GLY:HA3	1:6:919:A:H5'	269.74	0.56
26:D4:29:HIS:NE2	26:D4:34:ASN:HA	2.20	0.56
30:D8:15:VAL:HA	30:D8:28:VAL:HA	2.84	0.56
32:E0:59:GLY:O	32:E0:61:SER:N	4.09	0.56
33:E1:96:LYS:O	33:E1:98:VAL:N	2.31	0.56
39:L2:14:SER:OG	39:L2:15:ILE:N	2.36	0.56
39:L2:206:PRO:HG3	39:L2:213:GLY:HA3	1.86	0.56
45:L8:161:GLU:OE2	51:M5:26:ARG:NH2	2.37	0.56
36:1:2854:U:O3'	47:M0:160:PRO:HB3	2.05	0.56
49:M3:8:PRO:HD3	54:M8:164:ARG:HB3	2.81	0.56
53:M7:13:LYS:HB3	53:M7:152:GLU:HB2	1.86	0.56
66:O0:13:LYS:HD2	66:O0:100:ILE:HG23	1.88	0.56
67:O1:10:ARG:HH12	67:O1:44:MET:HG2	7.01	0.56
71:O5:85:THR:HG22	71:O5:87:ALA:N	2.20	0.56
2:S0:158:VAL:H	23:D1:69:LEU:HD12	1.71	0.56
3:S1:171:ILE:HA	3:S1:174:LYS:HE3	1.87	0.56
3:S1:36:SER:O	3:S1:38:PHE:N	2.38	0.56
36:1:343:U:O2	36:1:1439:U:H1'	2.04	0.56
1:2:1487:A:H2'	1:2:1488:G:C8	2.40	0.56
1:2:1482:C:OP2	1:2:1521:G:N2	2.36	0.56
38:4:155:A:OP1	45:L8:185:ARG:NH2	2.38	0.56
36:5:2403:G:H5'	36:5:2872:A:C2	2.41	0.56
26:D4:37:LYS:NZ	1:6:523:G:OP2	413.83	0.56
16:C4:121:VAL:O	1:6:886:U:O2'	287.78	0.56
1:6:973:A:H2'	1:6:974:A:C8	2.40	0.56
21:C9:57:ARG:HG2	21:C9:104:VAL:HG21	1.86	0.56
22:D0:51:VAL:HG13	22:D0:94:GLU:HB2	1.86	0.56
1:2:1796:C:OP2	28:D6:5:ARG:NH2	2.38	0.56
42:L5:269:SER:O	37:7:22:A:N6	323.24	0.56
45:L8:133:LYS:HB2	45:L8:199:ALA:O	3.91	0.56
46:L9:23:ARG:NH2	46:L9:41:ILE:O	5.96	0.56
46:L9:86:TYR:CE2	46:L9:151:VAL:HG22	2.42	0.56
58:N2:99:LYS:HG3	58:N2:102:GLU:HB2	1.88	0.56
36:1:1110:U:H2'	36:1:1111:U:C6	2.40	0.56
36:1:112:U:O2'	36:1:113:C:O5'	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:385:A:H2'	36:1:386:A:H8	1.70	0.56
1:2:304:U:H2'	1:2:305:C:H6	1.69	0.56
13:C1:18:HIS:O	13:C1:20:PHE:N	2.37	0.56
19:C7:47:ARG:NH1	19:C7:48:ASN:OD1	2.93	0.56
20:C8:97:ASP:N	20:C8:97:ASP:OD2	2.38	0.56
45:L8:134:TYR:H	45:L8:134:TYR:HD2	2.00	0.56
36:1:1447:G:H3'	53:M7:67:ILE:HD11	1.87	0.56
54:M8:23:ASN:HB3	54:M8:26:LEU:HB2	1.86	0.56
59:N3:11:PHE:CG	59:N3:88:ARG:HD2	3.14	0.56
79:Q3:81:SER:OG	79:Q3:82:THR:N	2.85	0.56
6:S4:163:ASP:HB2	6:S4:167:GLY:O	6.47	0.56
6:S4:65:LEU:HG	6:S4:70:VAL:HG11	1.87	0.56
7:S5:118:LEU:HD22	7:S5:129:PRO:HB2	1.88	0.56
7:S5:204:GLY:HA2	7:S5:211:ILE:HD12	2.72	0.56
36:1:1405:U:OP2	68:O2:59:SER:OG	2.23	0.56
36:1:107:A:H1'	36:1:325:A:N3	2.20	0.56
1:2:625:C:H2'	1:2:626:U:C6	2.41	0.56
36:5:1667:A:H2'	36:5:1668:G:C8	2.40	0.56
8:S6:149:LYS:NZ	1:6:141:U:OP1	313.30	0.56
14:C2:85:LYS:HG3	14:C2:87:PRO:HD3	3.17	0.56
5:S3:7:LYS:HE3	22:D0:27:THR:HG21	3.46	0.56
1:2:632:U:H4'	25:D3:11:SER:HB3	1.88	0.56
39:L2:225:ILE:HG21	39:L2:234:LYS:HA	1.88	0.56
46:L9:31:ARG:HB3	46:L9:149:ASN:HD21	3.20	0.56
48:M1:81:GLU:OE2	48:M1:89:TYR:OH	4.93	0.56
51:M5:201:ARG:NH2	36:5:692:A:OP1	96.86	0.56
59:N3:84:SER:HA	59:N3:94:TYR:HB3	2.04	0.56
67:O1:31:ARG:HB3	67:O1:31:ARG:NH1	4.86	0.56
67:O1:41:LYS:HE3	67:O1:45:GLY:HA2	1.88	0.56
36:1:181:U:H4'	73:O7:75:LYS:HD2	1.87	0.56
36:1:824:C:O2'	36:1:1534:A:N3	2.34	0.56
36:1:2603:G:H2'	36:1:2604:U:O4'	2.06	0.56
1:2:1175:U:H2'	1:2:1176:G:C8	2.41	0.56
1:2:1650:U:H2'	1:2:1651:A:C8	2.41	0.56
37:3:45:A:H2'	37:3:46:A:C8	2.41	0.56
41:L4:195:ARG:NH2	36:5:341:G:N7	109.77	0.56
36:5:58:G:O2'	36:5:61:A:H5'	2.06	0.56
21:C9:119:LYS:NZ	1:6:1369:U:OP1	441.55	0.56
37:7:4:U:H2'	37:7:5:G:C8	2.41	0.56
17:C5:102:PHE:HZ	1:6:1241:G:H5''	386.27	0.56
17:C5:81:ARG:NH1	17:C5:97:TYR:O	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:C6:47:LYS:HZ1	18:C6:114:ARG:HG2	1.71	0.56
18:C6:47:LYS:NZ	18:C6:114:ARG:HG2	2.21	0.56
19:C7:14:LYS:HG3	19:C7:69:ILE:HG12	5.90	0.56
21:C9:10:ALA:O	21:C9:13:ASP:N	2.35	0.56
21:C9:38:LYS:O	21:C9:39:THR:OG1	2.31	0.56
25:D3:31:LYS:HD2	25:D3:36:THR:HG21	1.88	0.56
16:C4:103:ARG:NH2	28:D6:48:ALA:O	6.45	0.56
39:L2:6:ARG:HH12	39:L2:199:THR:H	1.68	0.56
41:L4:185:LYS:HA	41:L4:201:GLN:HB3	3.01	0.56
41:L4:44:LYS:HB3	41:L4:47:ARG:HH11	1.71	0.56
46:L9:10:ILE:HD13	46:L9:75:VAL:HG11	1.87	0.56
56:N0:77:VAL:HG11	56:N0:106:LEU:HD11	1.87	0.56
6:S4:118:GLU:HA	6:S4:121:TYR:CE1	2.63	0.56
36:1:1352:A:H4'	36:1:1353:U:OP1	2.04	0.56
36:1:1654:A:H2'	36:1:1655:G:H5'	1.86	0.56
36:1:2694:A:C6	36:1:2695:A:C6	2.94	0.56
1:2:1042:G:H22	1:2:1076:A:H2	1.54	0.56
1:2:1655:A:N6	1:2:1745:G:O2'	2.38	0.56
36:5:1499:C:H42	36:5:1517:G:H1	1.53	0.56
15:C3:109:LYS:HD2	1:6:975:C:H5''	282.65	0.56
14:C2:54:ARG:NH1	14:C2:56:GLU:OE2	2.91	0.56
15:C3:119:GLU:HA	15:C3:122:ILE:HD12	1.86	0.56
21:C9:10:ALA:O	21:C9:12:GLN:N	2.39	0.56
28:D6:23:CYS:HB3	28:D6:28:LYS:H	2.10	0.56
40:L3:185:GLY:O	40:L3:191:LYS:NZ	3.75	0.56
41:L4:181:VAL:O	41:L4:182:LEU:HB2	2.05	0.56
44:L7:195:PHE:O	44:L7:199:ASN:HB3	2.80	0.56
46:L9:49:ASN:C	46:L9:51:GLN:H	2.09	0.56
51:M5:38:ARG:CZ	51:M5:60:VAL:HG13	2.36	0.56
57:N1:116:ARG:NH2	36:5:1097:G:N7	246.91	0.56
58:N2:59:ASP:OD1	58:N2:60:GLY:N	4.52	0.56
66:O0:26:GLY:O	66:O0:30:THR:HG23	2.49	0.56
71:O5:47:VAL:O	71:O5:51:ILE:HG13	2.06	0.56
3:S1:30:PHE:HB3	3:S1:96:LEU:HD22	3.63	0.56
9:S7:50:ASP:HA	9:S7:56:LYS:HA	1.87	0.56
11:S9:107:ARG:NH1	11:S9:112:GLN:OE1	2.38	0.56
1:2:1426:C:H5''	35:SM:93:ARG:HH12	1.69	0.56
36:1:1054:A:H5''	36:1:2637:A:N6	2.22	0.55
36:1:1349:G:H2'	36:1:1350:A:N3	2.21	0.55
36:1:1514:G:HO2'	36:1:1841:A:H2	1.54	0.55
36:1:3095:U:H2'	36:1:3096:C:C6	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:979:U:O2'	36:1:980:A:N7	2.28	0.55
1:2:1394:G:N2	1:2:1404:C:N3	2.43	0.55
36:5:2503:G:H1'	36:5:2504:U:H5	1.71	0.55
57:N1:17:ARG:HG2	36:5:2700:G:H5''	265.04	0.55
26:D4:53:ASP:HB3	26:D4:96:LEU:HD21	3.37	0.55
31:D9:19:ARG:HD2	31:D9:32:ARG:HD2	1.88	0.55
47:M0:87:LEU:HA	47:M0:138:VAL:HG13	1.88	0.55
50:M4:120:VAL:HG11	52:M6:199:TYR:CD2	2.62	0.55
36:1:1095:U:O2	57:N1:129:LYS:N	2.39	0.55
59:N3:10:LYS:NZ	59:N3:56:ASP:OD1	2.39	0.55
59:N3:38:ALA:HB3	59:N3:59:MET:HB2	1.87	0.55
60:N4:9:SER:OG	60:N4:10:GLY:N	2.37	0.55
4:S2:139:ILE:HG22	4:S2:141:ARG:HG2	2.57	0.55
5:S3:217:ILE:HG22	5:S3:218:LEU:H	1.70	0.55
10:S8:22:ARG:NE	10:S8:25:ARG:HD2	5.32	0.55
1:2:1160:A:H2'	1:2:1161:C:H6	1.70	0.55
5:S3:75:LYS:HB3	12:C0:22:VAL:HG22	2.40	0.55
13:C1:13:PHE:CE2	13:C1:15:LYS:HB3	2.41	0.55
24:D2:70:ASN:ND2	24:D2:130:TYR:O	2.30	0.55
24:D2:5:SER:O	24:D2:7:LEU:N	3.34	0.55
1:2:1795:U:H3'	28:D6:5:ARG:HH22	1.71	0.55
40:L3:187:SER:HB3	40:L3:190:GLU:HG3	1.89	0.55
40:L3:284:ARG:HH21	40:L3:293:ASN:C	3.54	0.55
41:L4:205:PRO:HG2	41:L4:225:VAL:HG22	3.86	0.55
44:L7:118:LYS:HG3	44:L7:191:VAL:HG11	1.87	0.55
42:L5:286:VAL:HG13	47:M0:206:LEU:HD22	1.89	0.55
49:M3:133:PRO:O	49:M3:135:ALA:N	3.23	0.55
36:1:3271:G:OP1	53:M7:171:ARG:HD3	2.06	0.55
54:M8:158:HIS:H	54:M8:186:VAL:HG22	4.71	0.55
50:M4:38:ILE:HD11	56:N0:150:PHE:HE2	1.69	0.55
56:N0:96:ASP:OD1	56:N0:97:VAL:N	2.36	0.55
63:N7:88:ASP:HB3	63:N7:121:ARG:HH12	1.70	0.55
69:O3:93:THR:O	69:O3:96:ALA:N	2.95	0.55
3:S1:59:ASP:HA	3:S1:62:LYS:HG3	1.88	0.55
6:S4:175:PHE:HE1	6:S4:225:VAL:HG11	2.40	0.55
8:S6:174:LYS:HG3	1:6:79:C:H1'	341.52	0.55
9:S7:69:GLY:HA2	9:S7:72:LYS:HD2	1.87	0.55
36:1:1301:A:H4'	36:1:1302:A:O5'	2.06	0.55
36:1:767:U:H1'	36:1:768:C:C6	2.42	0.55
36:1:97:U:O4	49:M3:11:LYS:NZ	2.29	0.55
1:2:1169:G:N1	1:2:1575:G:OP2	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1335:U:H2'	1:2:1336:A:H8	1.71	0.55
1:2:1564:U:H2'	1:2:1565:C:C6	2.41	0.55
1:2:649:U:O2'	1:2:650:U:O4'	2.24	0.55
57:N1:129:LYS:HB2	36:5:1098:A:O5'	252.38	0.55
36:5:1472:U:H2'	36:5:1473:G:H8	1.71	0.55
36:5:1838:G:H5''	36:5:1839:A:H5'	1.89	0.55
36:5:2261:G:O2'	36:5:2263:C:N4	2.39	0.55
1:6:1648:A:H2'	1:6:1649:G:C8	2.40	0.55
18:C6:77:GLN:O	18:C6:81:ILE:HG13	4.05	0.55
22:D0:118:VAL:HG13	22:D0:119:ALA:H	1.70	0.55
29:D7:36:LYS:HE2	29:D7:78:SER:HB3	1.89	0.55
40:L3:303:LYS:HD2	40:L3:361:THR:HG21	1.90	0.55
48:M1:7:ASN:N	48:M1:8:PRO:HD3	2.68	0.55
51:M5:186:GLY:O	51:M5:190:THR:HG23	2.34	0.55
66:O0:34:LEU:HD23	66:O0:59:TYR:HB3	1.88	0.55
2:S0:118:PRO:HG2	2:S0:141:ILE:HD13	1.92	0.55
36:1:3029:A:C5	36:1:3030:G:H1'	2.41	0.55
36:1:3276:G:H5'	43:L6:48:ARG:NH2	2.21	0.55
1:2:1147:A:H2'	1:2:1148:C:C6	2.42	0.55
41:L4:193:LYS:NZ	36:5:1420:C:OP2	111.81	0.55
36:5:2157:G:N1	36:5:2178:A:OP2	2.32	0.55
53:M7:139:TYR:CE2	36:5:2355:G:H4'	147.46	0.55
36:5:174:C:H42	36:5:244:G:H1	1.53	0.55
36:5:2451:G:N2	36:5:2496:C:O2	2.39	0.55
1:6:1272:U:H3	1:6:1438:G:H1	1.54	0.55
1:6:217:A:O2'	1:6:218:A:O5'	2.20	0.55
1:6:846:G:H2'	1:6:847:A:C8	2.41	0.55
15:C3:15:ALA:O	1:6:959:U:H5''	350.36	0.55
37:7:91:G:H2'	37:7:92:A:C8	2.42	0.55
30:D8:7:VAL:HG13	30:D8:55:VAL:HG13	1.87	0.55
39:L2:116:VAL:HG22	39:L2:126:LEU:HB2	3.32	0.55
40:L3:238:LEU:HB3	40:L3:242:THR:HG21	2.16	0.55
40:L3:70:ARG:HB3	40:L3:70:ARG:HH11	4.32	0.55
41:L4:138:ARG:NH1	41:L4:138:ARG:O	3.01	0.55
51:M5:185:ALA:HB3	51:M5:190:THR:HG22	2.11	0.55
56:N0:117:ARG:NH2	36:5:1322:U:OP1	280.93	0.55
59:N3:28:ASN:HD21	59:N3:112:SER:H	2.14	0.55
60:N4:3:VAL:HG11	60:N4:12:LYS:HD3	2.09	0.55
49:M3:49:ARG:NH2	71:O5:113:GLN:OE1	4.49	0.55
2:S0:50:VAL:H	19:C7:109:LEU:HD21	1.92	0.55
7:S5:77:TYR:CZ	7:S5:87:CYS:HB2	2.59	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:S8:141:ARG:NH2	1:6:196:G:O6	277.48	0.55
11:S9:63:ASP:OD1	11:S9:64:GLU:N	3.47	0.55
36:1:1665:C:H2'	36:1:1666:G:C8	2.40	0.55
35:SM:33:LYS:HE2	36:5:2667:A:H5''	285.97	0.55
50:M4:106:ARG:NH2	36:5:3207:U:OP1	302.34	0.55
1:6:1068:C:H2'	1:6:1069:A:H8	1.71	0.55
1:6:793:A:H5''	1:6:794:U:C5	2.41	0.55
17:C5:122:THR:HG21	1:6:1455:G:OP1	370.08	0.55
28:D6:23:CYS:HB2	28:D6:74:CYS:SG	4.29	0.55
40:L3:92:TYR:O	40:L3:156:SER:N	2.34	0.55
42:L5:95:TRP:CH2	42:L5:181:PRO:HD3	4.77	0.55
42:L5:183:TRP:CH2	42:L5:188:GLU:HA	2.41	0.55
49:M3:83:ALA:HA	49:M3:117:LYS:HE3	1.89	0.55
57:N1:88:ARG:HB2	36:5:2722:U:H4'	216.49	0.55
61:N5:105:VAL:HG12	61:N5:106:ASP:O	2.07	0.55
63:N7:51:LEU:HB3	63:N7:65:ARG:HH11	2.48	0.55
9:S7:51:VAL:HG23	9:S7:53:GLY:H	1.72	0.55
1:2:1677:C:OP1	10:S8:42:ARG:NH1	2.39	0.55
34:SR:200:ASN:H	34:SR:215:GLY:HA2	2.44	0.55
34:SR:248:ASN:OD1	34:SR:249:ARG:N	4.41	0.55
36:1:3159:C:H2'	36:1:3160:U:C6	2.42	0.55
36:1:3259:U:H6	36:1:3259:U:H5'	1.72	0.55
1:2:1342:C:OP1	34:SR:102:ARG:NH2	2.39	0.55
1:2:1696:G:N2	1:2:1705:C:H41	2.05	0.55
37:3:92:A:C5	37:3:93:C:H1'	2.41	0.55
63:N7:48:ARG:NH2	36:5:1631:C:OP2	192.33	0.55
36:5:2103:U:H2'	36:5:2104:A:H8	1.72	0.55
1:6:1045:C:H2'	1:6:1046:G:H8	1.71	0.55
1:6:65:A:H2	1:6:84:A:H62	1.55	0.55
14:C2:80:ASN:HA	14:C2:86:VAL:HG12	1.87	0.55
27:D5:43:ASP:O	27:D5:45:GLU:N	2.40	0.55
31:D9:14:TYR:OH	1:6:1553:G:O2'	403.17	0.55
40:L3:4:ARG:HD2	40:L3:7:GLU:HG3	1.87	0.55
46:L9:4:ILE:HD11	56:N0:150:PHE:CD2	2.41	0.55
61:N5:91:ASN:O	61:N5:93:TYR:N	3.18	0.55
62:N6:17:LYS:O	62:N6:21:THR:OG1	2.18	0.55
36:1:1191:U:H3'	76:Q0:113:ARG:NH2	2.21	0.55
7:S5:151:GLY:HA3	7:S5:155:ALA:HA	4.67	0.55
11:S9:34:PHE:HD1	11:S9:111:THR:HG21	1.94	0.55
36:1:210:U:C2	36:1:230:U:H4'	2.42	0.55
36:1:3326:G:H2'	36:1:3327:G:H8	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1318:G:H2'	1:2:1319:A:C8	2.42	0.55
1:2:87:C:O2'	1:2:169:A:N1	2.30	0.55
36:5:1015:U:O3'	36:5:1016:C:H2'	2.07	0.55
36:5:1047:A:N3	36:5:2633:U:O2'	2.37	0.55
36:5:900:G:H1'	36:5:1589:A:H61	1.72	0.55
36:5:2619:G:H1	84:B:75:CH:HN3	1.55	0.55
1:6:152:U:C2	1:6:163:G:N2	2.75	0.55
1:6:75:U:O2'	1:6:76:A:O5'	2.23	0.55
15:C3:94:LYS:HG2	15:C3:118:ILE:HD13	2.01	0.55
5:S3:211:PRO:HG3	19:C7:20:TYR:CZ	2.41	0.55
19:C7:57:LEU:HA	19:C7:60:ARG:HG2	1.88	0.55
20:C8:88:ARG:NH1	20:C8:112:ASP:OD2	2.73	0.55
21:C9:117:SER:HB3	21:C9:123:ARG:HB2	1.88	0.55
25:D3:134:ALA:HB1	25:D3:140:LYS:HG3	3.15	0.55
1:2:778:G:H22	26:D4:10:ARG:NH1	2.04	0.55
33:E1:144:CYS:O	33:E1:146:SER:N	2.39	0.55
47:M0:48:LEU:HD22	47:M0:49:CYS:H	1.72	0.55
48:M1:81:GLU:HB2	48:M1:167:TYR:HE2	5.04	0.55
52:M6:10:ASP:CG	52:M6:37:ARG:HH21	2.34	0.55
49:M3:159:VAL:HA	64:N8:124:ILE:HD11	2.25	0.55
64:N8:19:LYS:HD2	64:N8:25:HIS:CD2	4.39	0.55
71:O5:14:LYS:NZ	71:O5:62:GLN:OE1	7.30	0.55
7:S5:79:ASN:OD1	7:S5:79:ASN:N	2.40	0.55
10:S8:36:THR:OG1	10:S8:96:LEU:O	2.24	0.55
36:1:1488:G:C2	36:1:1489:A:C8	2.95	0.55
36:1:1683:A:H2'	36:1:1684:U:O4'	2.07	0.55
36:1:27:C:O2'	36:1:327:A:N3	2.34	0.55
36:1:438:A:O2'	36:1:495:G:O2'	2.11	0.55
36:1:916:G:H5'	36:1:917:A:OP1	2.07	0.55
1:2:1018:U:H2'	1:2:1019:A:C8	2.40	0.55
36:5:1657:C:O2'	36:5:1797:A:OP2	2.14	0.55
36:5:2541:U:H4'	36:5:2542:U:OP1	2.07	0.55
24:D2:2:THR:N	1:6:1034:C:HO2'	337.83	0.55
1:6:895:G:H22	1:6:917:U:H3	1.54	0.55
17:C5:85:ILE:HG22	17:C5:112:LEU:HD23	2.22	0.55
20:C8:94:ASP:OD2	20:C8:95:GLY:N	2.40	0.55
22:D0:37:VAL:HG13	22:D0:107:THR:HG22	4.27	0.55
24:D2:37:PHE:CD2	24:D2:103:ILE:HD12	2.42	0.55
1:2:571:G:H5''	25:D3:114:LYS:HD3	1.88	0.55
49:M3:92:THR:HG21	71:O5:111:PHE:HB3	3.53	0.55
44:L7:69:ALA:HB2	57:N1:140:ILE:HD13	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:O7:31:LYS:O	73:O7:33:THR:HG22	2.07	0.55
2:S0:172:LEU:O	2:S0:176:LEU:HG	2.07	0.55
8:S6:18:ILE:O	8:S6:20:ASP:N	3.48	0.55
36:1:1277:C:HO2'	36:1:1278:A:H8	1.54	0.55
36:1:1397:C:H2'	36:1:1398:U:H6	1.72	0.55
36:1:1427:U:OP2	64:N8:4:ARG:NH2	2.36	0.55
36:1:1590:G:N2	36:1:1798:A:N1	2.53	0.55
36:1:282:G:C8	36:1:282:G:H3'	2.42	0.55
36:1:964:G:O2'	64:N8:41:HIS:NE2	2.36	0.55
1:2:435:C:H2'	1:2:436:A:C8	2.42	0.55
1:2:652:G:N2	1:2:682:C:N3	2.46	0.55
64:N8:4:ARG:NH2	36:5:1427:U:OP2	135.49	0.55
21:C9:68:ARG:NH1	1:6:1521:G:O6	413.33	0.55
1:6:235:G:H2'	1:6:236:A:C8	2.40	0.55
38:8:157:U:H2'	38:8:158:U:C6	2.42	0.55
14:C2:62:LEU:HD22	14:C2:75:VAL:HG11	1.89	0.55
18:C6:82:ARG:HH22	18:C6:114:ARG:HG3	4.02	0.55
1:2:1548:G:H5'	20:C8:89:GLN:O	2.07	0.55
22:D0:80:GLU:OE2	31:D9:54:LYS:NZ	4.45	0.55
49:M3:47:ALA:HB3	49:M3:49:ARG:HG3	4.57	0.55
51:M5:96:ARG:NH2	51:M5:100:ALA:HB1	2.40	0.55
36:1:269:G:H5'	51:M5:120:TRP:CE3	2.41	0.55
63:N7:124:ALA:O	63:N7:126:LYS:N	2.37	0.55
68:O2:102:ALA:HB2	68:O2:125:ARG:HG2	4.29	0.55
68:O2:40:SER:O	68:O2:44:ARG:HG3	2.06	0.55
36:1:351:A:N6	75:O9:37:TYR:O	2.37	0.55
6:S4:184:THR:OG1	6:S4:224:ASN:O	3.04	0.55
8:S6:126:ASP:OD2	8:S6:127:THR:N	2.70	0.55
36:1:263:C:H2'	36:1:264:G:O4'	2.07	0.55
1:2:108:A:H2'	1:2:109:G:C8	2.42	0.55
1:2:1241:G:H5'	17:C5:102:PHE:CZ	2.41	0.55
1:2:138:A:OP2	1:2:1706:C:O2'	2.17	0.55
1:2:1681:A:H2'	1:2:1682:U:H5'	1.89	0.55
1:6:1175:U:H2'	1:6:1176:G:C8	2.42	0.55
1:6:1688:U:H3	1:6:1713:G:H1	1.55	0.55
19:C7:20:TYR:CE1	19:C7:38:ILE:HD11	2.42	0.55
22:D0:99:ILE:HD12	22:D0:102:ARG:HD3	1.89	0.55
23:D1:17:CYS:HB2	23:D1:56:SER:HB3	2.02	0.55
1:2:1104:U:C5	25:D3:4:GLY:HA2	2.42	0.55
27:D5:92:ILE:HD11	27:D5:100:ILE:HG22	2.54	0.55
28:D6:79:ILE:HA	28:D6:84:VAL:HG11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:D8:42:ARG:NH2	30:D8:58:GLU:HG3	2.21	0.55
36:1:3037:U:H5''	40:L3:348:ARG:NH1	2.22	0.55
42:L5:277:LEU:HD12	42:L5:282:ARG:HG3	5.49	0.55
46:L9:18:VAL:HB	46:L9:27:VAL:HG22	3.57	0.55
50:M4:40:ASP:OD1	50:M4:42:LYS:N	2.36	0.55
52:M6:157:GLU:O	52:M6:161:LYS:N	2.25	0.55
62:N6:119:ILE:HG22	62:N6:124:GLY:HA3	2.46	0.55
66:O0:16:LEU:HD13	66:O0:19:LYS:HE3	1.89	0.55
4:S2:154:LEU:HD12	4:S2:155:ALA:H	1.71	0.55
36:1:1238:C:N4	36:1:1245:A:OP2	2.40	0.54
36:1:915:A:H8	36:1:2136:C:O2'	1.90	0.54
36:1:3258:U:O2'	36:1:3260:G:OP1	2.17	0.54
36:1:507:U:H2'	36:1:508:U:C6	2.42	0.54
1:2:327:U:H2'	1:2:328:A:C8	2.43	0.54
36:5:2836:C:H5	36:5:2852:C:N4	2.03	0.54
49:M3:7:LEU:HD22	36:5:668:G:H1'	159.40	0.54
17:C5:126:VAL:HG13	17:C5:127:ARG:H	1.71	0.54
22:D0:109:GLU:HB3	22:D0:112:VAL:HB	1.88	0.54
40:L3:41:VAL:HA	40:L3:185:GLY:CA	2.37	0.54
48:M1:16:LYS:H	48:M1:130:VAL:HG13	1.72	0.54
49:M3:140:SER:OG	49:M3:141:ALA:N	2.40	0.54
52:M6:22:VAL:O	52:M6:26:GLN:HG2	2.07	0.54
36:1:744:A:H4'	54:M8:142:GLY:O	2.08	0.54
54:M8:176:ARG:NH1	64:N8:46:ASP:OD2	2.52	0.54
54:M8:30:VAL:O	54:M8:34:THR:HG23	2.37	0.54
57:N1:86:GLU:OE1	57:N1:88:ARG:NH1	2.40	0.54
63:N7:128:GLN:O	63:N7:130:PHE:N	2.90	0.54
2:S0:70:PRO:O	2:S0:95:ALA:N	2.30	0.54
4:S2:214:ALA:O	4:S2:218:ILE:HG13	2.22	0.54
34:SR:19:TRP:HB2	34:SR:38:ARG:HD2	1.92	0.54
34:SR:267:PRO:HD2	34:SR:269:TYR:HE1	2.85	0.54
36:1:1582:C:O2'	36:1:1583:A:O5'	2.21	0.54
36:1:2150:G:O2'	36:1:2189:U:OP1	2.19	0.54
36:1:2767:U:H2'	36:1:2768:U:C6	2.42	0.54
36:1:3117:C:H2'	36:1:3118:C:O4'	2.06	0.54
1:2:1036:A:H1'	24:D2:9:ASP:OD1	2.08	0.54
1:2:45:U:O2'	1:2:46:A:H2'	2.07	0.54
1:6:206:A:H1'	1:6:262:U:C2	2.42	0.54
1:6:846:G:H2'	1:6:847:A:H8	1.72	0.54
37:7:4:U:H2'	37:7:5:G:H8	1.72	0.54
17:C5:18:ARG:HG2	20:C8:92:ILE:HA	2.27	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:C7:27:ASP:OD2	19:C7:30:THR:N	3.47	0.54
37:3:5:G:O3'	42:L5:54:ARG:HG3	2.07	0.54
44:L7:145:ARG:HA	44:L7:185:ILE:HD13	1.89	0.54
47:M0:210:ILE:HA	47:M0:217:PHE:CE2	2.66	0.54
51:M5:102:ALA:O	51:M5:106:VAL:HG13	2.13	0.54
54:M8:158:HIS:H	54:M8:186:VAL:CG1	2.20	0.54
62:N6:69:LYS:O	62:N6:83:ASP:N	3.44	0.54
70:O4:37:LYS:HB2	70:O4:58:ARG:HH21	1.72	0.54
4:S2:227:PRO:HA	4:S2:230:TRP:CD1	2.79	0.54
36:1:1240:A:H3'	36:1:1241:U:H5'	1.89	0.54
36:1:1724:U:H1'	36:1:1725:C:C6	2.43	0.54
36:1:2378:C:H2'	36:1:2379:U:C6	2.42	0.54
36:1:2812:C:H2'	36:1:2813:A:H8	1.72	0.54
1:2:427:C:O2'	1:2:459:G:N3	2.31	0.54
36:5:230:U:H2'	36:5:231:G:O4'	2.08	0.54
36:5:2378:C:H2'	36:5:2379:U:H6	1.72	0.54
57:N1:101:CYS:HB3	36:5:990:U:H1'	251.31	0.54
18:C6:139:GLN:NE2	1:6:1465:C:OP1	352.71	0.54
42:L5:152:ARG:HG3	37:7:44:C:H4'	282.86	0.54
13:C1:109:VAL:HG11	13:C1:125:VAL:HG21	3.94	0.54
40:L3:346:THR:O	40:L3:348:ARG:N	3.95	0.54
41:L4:264:SER:OG	41:L4:267:VAL:N	4.36	0.54
41:L4:74:ILE:HD12	41:L4:75:PRO:HD2	5.03	0.54
42:L5:244:HIS:HA	42:L5:247:ILE:HD12	1.89	0.54
44:L7:229:PHE:CD1	44:L7:229:PHE:C	2.97	0.54
45:L8:84:ARG:HH22	45:L8:181:LYS:NZ	2.05	0.54
46:L9:104:VAL:HG22	46:L9:113:GLU:HB2	1.88	0.54
47:M0:144:ASN:O	47:M0:146:ASP:N	2.90	0.54
47:M0:71:CYS:SG	47:M0:72:ALA:N	3.49	0.54
49:M3:28:GLN:HB3	51:M5:201:ARG:HD3	1.87	0.54
51:M5:47:LYS:O	51:M5:50:ARG:HG2	2.07	0.54
54:M8:170:ARG:O	54:M8:171:LYS:HB2	4.08	0.54
37:3:97:A:OP1	56:N0:40:ARG:NH1	2.41	0.54
57:N1:78:LYS:N	57:N1:85:LEU:O	3.50	0.54
64:N8:28:HIS:CD2	64:N8:32:ARG:HG2	2.92	0.54
67:O1:49:VAL:HG22	67:O1:91:SER:HB2	1.87	0.54
6:S4:125:LYS:O	6:S4:142:HIS:N	2.71	0.54
7:S5:42:LEU:HB2	7:S5:46:TRP:O	2.07	0.54
9:S7:74:GLN:NE2	9:S7:92:PHE:HB2	2.22	0.54
10:S8:197:THR:HA	10:S8:200:LYS:HZ3	1.71	0.54
11:S9:149:ARG:O	11:S9:151:ASP:N	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1103:A:H4'	36:1:1103:A:OP2	2.07	0.54
36:1:150:A:OP1	51:M5:56:LYS:NZ	2.35	0.54
36:1:2102:U:H2'	36:1:2103:U:C6	2.43	0.54
36:1:274:G:H2'	36:1:275:U:O4'	2.07	0.54
36:1:309:U:OP1	72:O6:84:LYS:NZ	2.38	0.54
1:2:1769:U:H1'	16:C4:136:ARG:HG3	1.88	0.54
1:2:732:G:O2'	1:2:733:A:O4'	2.25	0.54
37:3:26:C:H5''	42:L5:56:THR:HB	1.89	0.54
36:5:1239:C:H42	36:5:1249:G:H22	1.53	0.54
36:5:3161:C:H2'	36:5:3162:C:C6	2.42	0.54
36:5:3276:G:OP2	36:5:3276:G:H2'	2.07	0.54
36:5:811:U:H2'	36:5:812:G:C8	2.42	0.54
1:6:619:A:N3	1:6:1141:G:H1'	2.23	0.54
1:6:27:U:H2'	1:6:28:A:C8	2.42	0.54
2:S0:52:LYS:HD2	23:D1:82:VAL:HA	1.89	0.54
26:D4:11:LYS:HB2	26:D4:24:VAL:HG23	2.01	0.54
32:E0:46:ASN:HD21	32:E0:48:THR:HG23	4.79	0.54
41:L4:170:LYS:HD3	41:L4:175:HIS:ND1	2.22	0.54
42:L5:153:THR:HG22	42:L5:179:ARG:HD2	1.89	0.54
44:L7:196:LYS:NZ	36:5:1101:G:OP2	241.36	0.54
45:L8:152:LEU:HB2	45:L8:198:ALA:HB3	1.89	0.54
45:L8:95:ASN:HA	45:L8:98:ARG:NH1	5.78	0.54
46:L9:70:THR:HB	36:5:3112:G:O2'	329.21	0.54
36:1:2674:A:H5''	48:M1:105:GLY:HA3	1.90	0.54
53:M7:141:SER:O	53:M7:143:PRO:HD3	2.51	0.54
54:M8:100:THR:HG22	54:M8:120:GLU:HB3	3.33	0.54
66:O0:20:SER:OG	66:O0:21:GLY:N	2.40	0.54
64:N8:14:HIS:CE1	68:O2:36:LYS:HE2	2.48	0.54
5:S3:102:ALA:HB1	5:S3:173:ARG:HG3	2.80	0.54
6:S4:206:ASP:HB2	6:S4:222:LEU:HB2	1.89	0.54
6:S4:230:GLU:HB2	6:S4:233:LYS:NZ	4.87	0.54
6:S4:31:PRO:HB2	6:S4:38:LEU:HB2	4.79	0.54
9:S7:86:GLN:HG2	9:S7:87:ASP:H	1.72	0.54
11:S9:102:GLU:HA	11:S9:105:LEU:HD23	4.37	0.54
11:S9:40:LYS:HE2	1:6:593:U:H5''	401.45	0.54
1:2:763:G:OP2	11:S9:79:ARG:NH1	2.38	0.54
34:SR:105:GLY:HA3	34:SR:134:TRP:HH2	2.93	0.54
36:1:2314:U:O2'	36:1:2315:G:OP1	2.20	0.54
36:1:860:G:C5	39:L2:181:LYS:HB2	2.41	0.54
1:2:755:A:C2	6:S4:13:ALA:HA	2.42	0.54
1:2:803:A:C4	9:S7:104:ARG:HG3	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:M0:119:TRP:HZ3	36:5:1126:G:H5''	256.29	0.54
36:5:1692:U:O4	36:5:1693:C:N4	2.41	0.54
36:5:2255:A:OP2	36:5:2261:G:N1	2.30	0.54
36:5:503:C:H2'	36:5:504:A:H8	1.72	0.54
19:C7:33:ARG:O	19:C7:37:GLU:HG2	2.08	0.54
55:M9:138:LEU:O	55:M9:142:ILE:HG13	2.32	0.54
55:M9:84:THR:O	55:M9:88:ARG:HG2	3.89	0.54
66:O0:99:ASP:O	66:O0:103:THR:OG1	3.04	0.54
68:O2:105:ARG:O	68:O2:109:LEU:HB2	2.70	0.54
69:O3:52:VAL:HG22	69:O3:66:VAL:HG22	1.88	0.54
71:O5:43:LYS:O	71:O5:46:THR:HG22	2.08	0.54
71:O5:64:GLU:O	71:O5:68:GLN:N	3.45	0.54
2:S0:139:VAL:HA	4:S2:62:PRO:HG3	1.90	0.54
4:S2:106:ASP:OD2	4:S2:110:HIS:ND1	2.40	0.54
8:S6:84:TYR:OH	8:S6:91:GLU:OE1	2.24	0.54
36:1:1097:G:H4'	36:1:1098:A:O5'	2.08	0.54
36:1:2601:A:H2'	36:1:2602:G:H8	1.73	0.54
1:2:158:U:O2'	1:2:159:U:H3'	2.07	0.54
1:2:711:U:H1'	1:2:712:G:H5'	1.90	0.54
36:5:1497:C:H2'	36:5:1498:A:H8	1.71	0.54
36:5:2773:C:H2'	36:5:2774:C:C6	2.43	0.54
36:5:411:U:H2'	36:5:412:G:H8	1.73	0.54
1:6:1620:C:H2'	1:6:1621:U:C6	2.42	0.54
1:6:25:C:OP2	1:6:25:C:H4'	2.06	0.54
19:C7:23:LYS:HB3	19:C7:34:LEU:HD11	1.90	0.54
20:C8:99:HIS:HD2	20:C8:101:LEU:HD21	1.92	0.54
21:C9:33:TYR:HD1	21:C9:34:VAL:H	3.62	0.54
23:D1:2:GLU:HG3	23:D1:6:GLY:HA2	5.66	0.54
41:L4:264:SER:OG	41:L4:265:GLU:N	2.40	0.54
53:M7:67:ILE:HG13	53:M7:82:ARG:CZ	2.38	0.54
66:O0:84:LEU:HD22	36:5:1715:A:N7	260.77	0.54
6:S4:17:HIS:HB2	6:S4:108:ARG:HA	1.90	0.54
34:SR:17:ASN:ND2	34:SR:17:ASN:O	2.40	0.54
36:1:2443:A:O2'	36:1:2444:C:H5'	2.07	0.54
36:1:2946:A:H5''	36:1:2947:G:H5'	1.89	0.54
36:1:59:G:H2'	38:4:33:A:O2'	2.06	0.54
1:2:1649:G:H2'	1:2:1650:U:C6	2.43	0.54
36:5:1528:G:H1	36:5:1832:C:H42	1.54	0.54
20:C8:63:GLN:O	20:C8:67:GLU:HB2	3.28	0.54
21:C9:14:PHE:HZ	21:C9:132:LEU:HD22	3.79	0.54
23:D1:71:ARG:HG3	23:D1:83:TRP:CE2	4.09	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:E1:135:HIS:HB2	33:E1:138:ARG:HG3	1.90	0.54
40:L3:230:THR:HA	40:L3:235:THR:HG22	1.89	0.54
40:L3:280:HIS:HB3	40:L3:324:VAL:HG13	1.89	0.54
42:L5:290:ILE:HG12	47:M0:206:LEU:HD11	5.00	0.54
53:M7:132:ALA:O	53:M7:135:ARG:NH2	3.25	0.54
53:M7:26:PHE:CE1	53:M7:120:ASN:HA	2.54	0.54
59:N3:45:ARG:HB3	59:N3:48:ARG:HG3	1.88	0.54
61:N5:105:VAL:HA	61:N5:130:TYR:CE2	2.55	0.54
69:O3:35:VAL:HG13	69:O3:40:ASP:HB3	1.88	0.54
9:S7:14:THR:OG1	9:S7:15:GLU:N	2.44	0.54
10:S8:10:LYS:NZ	1:6:337:G:O2'	284.48	0.54
36:1:1362:G:H2'	36:1:1363:A:C8	2.42	0.54
36:1:2093:A:H3'	36:1:2093:A:N3	2.22	0.54
36:1:348:A:H1'	36:1:352:A:O2'	2.08	0.54
36:1:385:A:H2'	36:1:386:A:C8	2.43	0.54
36:1:426:G:OP1	68:O2:15:LYS:NZ	2.39	0.54
1:2:1682:U:O2'	1:2:1683:C:H5'	2.06	0.54
1:2:233:C:H4'	1:2:234:G:C2	2.42	0.54
1:2:589:C:H2'	1:2:590:C:H6	1.72	0.54
1:2:372:G:N2	1:2:612:U:O4	2.29	0.54
1:2:826:U:H2'	1:2:827:C:C6	2.43	0.54
36:5:761:A:C2	36:5:771:A:H1'	2.43	0.54
1:6:729:G:O2'	1:6:730:G:O5'	2.25	0.54
61:N5:56:ARG:HG2	38:8:134:G:OP1	78.90	0.54
18:C6:130:GLY:HA3	18:C6:137:ARG:HH12	2.86	0.54
28:D6:43:ASN:OD1	28:D6:66:LYS:NZ	6.83	0.54
28:D6:87:ARG:HB3	28:D6:91:ASP:HB3	1.89	0.54
41:L4:44:LYS:HB3	41:L4:47:ARG:NH1	2.30	0.54
42:L5:155:THR:HA	42:L5:179:ARG:HD3	2.14	0.54
45:L8:54:GLU:O	45:L8:58:VAL:HG23	2.32	0.54
46:L9:73:SER:OG	46:L9:74:LEU:N	2.39	0.54
56:N0:77:VAL:HG13	56:N0:126:VAL:HG22	1.97	0.54
58:N2:32:SER:HA	58:N2:35:LYS:HB3	1.89	0.54
64:N8:47:LYS:C	64:N8:49:HIS:H	2.25	0.54
69:O3:51:TYR:HB3	69:O3:67:MET:HB2	2.53	0.54
38:4:53:A:C2	75:O9:35:ILE:HD11	2.43	0.54
4:S2:116:LYS:HG2	4:S2:127:ALA:HB3	1.90	0.54
8:S6:31:ARG:N	8:S6:34:GLN:OE1	2.36	0.54
1:2:768:C:C2	11:S9:143:ILE:HD13	2.43	0.54
34:SR:70:ASP:HB3	34:SR:113:VAL:HG12	2.26	0.54
1:2:1487:A:OP1	31:D9:34:TYR:OH	2.14	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:472:U:H5''	11:S9:11:THR:HG23	1.90	0.54
1:2:505:A:H61	1:2:507:U:H3	1.55	0.54
36:5:2651:G:C4	36:5:2796:G:C2	2.96	0.54
36:5:2712:U:H2'	36:5:2713:U:H6	1.73	0.54
1:6:1079:U:H2'	1:6:1080:U:C6	2.41	0.54
18:C6:135:ARG:NH1	1:6:1583:A:OP1	382.61	0.54
1:6:654:C:H2'	1:6:655:G:C8	2.43	0.54
1:6:918:U:H2'	1:6:919:A:C8	2.38	0.54
14:C2:132:GLU:HG2	14:C2:133:LEU:HD12	3.04	0.54
14:C2:68:GLU:O	14:C2:70:ASN:N	2.40	0.54
14:C2:70:ASN:HA	14:C2:73:LYS:HE2	2.12	0.54
18:C6:42:GLU:HA	18:C6:45:ARG:HB2	1.89	0.54
26:D4:25:VAL:N	26:D4:71:GLY:O	2.93	0.54
43:L6:3:ALA:HB1	68:O2:75:LEU:HD13	3.09	0.54
45:L8:195:SER:O	45:L8:197:VAL:N	3.33	0.54
45:L8:75:ILE:HG22	45:L8:76:ALA:H	1.73	0.54
46:L9:109:ALA:O	46:L9:110:LYS:HB2	2.26	0.54
47:M0:5:PRO:C	47:M0:7:ARG:H	2.10	0.54
51:M5:123:GLN:HB2	51:M5:128:LYS:HG2	2.64	0.54
63:N7:54:THR:OG1	63:N7:55:LYS:N	2.40	0.54
51:M5:147:ARG:NH1	71:O5:101:THR:OG1	4.01	0.54
71:O5:7:TYR:HA	71:O5:10:ARG:HG3	3.63	0.54
72:O6:97:SER:HB3	72:O6:98:ARG:HE	1.73	0.54
6:S4:131:LEU:HD12	1:6:252:U:H4'	324.81	0.54
36:1:2565:U:H2'	36:1:2566:C:C6	2.42	0.54
36:1:715:A:H8	64:N8:115:LYS:HG2	1.73	0.54
1:2:1002:G:N2	1:2:1760:G:O3'	2.41	0.54
1:2:1761:U:O2'	1:2:1762:A:OP2	2.24	0.54
63:N7:73:LYS:HE2	36:5:1636:U:H5''	211.82	0.54
66:O0:28:LYS:NZ	36:5:1713:G:O6	236.16	0.54
13:C1:83:THR:HG21	1:6:325:G:H4'	288.37	0.54
36:1:3147:G:O2'	40:L3:104:THR:OG1	2.20	0.54
40:L3:114:VAL:HG22	40:L3:163:HIS:CE1	2.46	0.54
43:L6:77:ARG:NH1	43:L6:78:ARG:O	2.40	0.54
45:L8:139:VAL:O	45:L8:143:ILE:HG13	2.26	0.54
45:L8:203:VAL:HG13	45:L8:207:ASP:HB2	1.90	0.54
46:L9:137:SER:HB3	46:L9:143:GLU:HB3	1.90	0.54
51:M5:68:ARG:HG3	36:5:291:C:OP1	145.73	0.54
58:N2:50:LEU:H	58:N2:50:LEU:HD12	1.72	0.54
78:Q2:17:CYS:SG	78:Q2:77:CYS:HB3	2.48	0.54
2:S0:69:ASN:HB3	2:S0:71:GLU:CD	2.28	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S0:70:PRO:HB2	2:S0:94:GLY:HA3	1.90	0.54
3:S1:126:THR:HG22	3:S1:136:ARG:HE	1.72	0.54
5:S3:99:VAL:HG22	5:S3:171:ALA:HB2	3.31	0.54
5:S3:66:ILE:O	5:S3:70:THR:N	2.90	0.54
6:S4:207:LEU:HD22	6:S4:221:ARG:HA	3.71	0.54
9:S7:50:ASP:HB3	9:S7:56:LYS:HG2	1.88	0.54
10:S8:197:THR:HG22	10:S8:200:LYS:HZ3	1.73	0.54
34:SR:160:GLU:O	34:SR:162:ALA:N	2.38	0.54
1:2:998:A:N1	1:2:1006:C:N4	2.57	0.53
37:3:60:G:H2'	37:3:61:G:H8	1.72	0.53
38:4:24:G:OP2	62:N6:13:ARG:NH1	2.29	0.53
36:5:1317:A:O2'	36:5:1318:A:H3'	2.08	0.53
44:L7:206:LYS:HB3	36:5:1334:U:H5''	236.41	0.53
36:5:1913:A:N3	36:5:2120:A:H2'	2.23	0.53
24:D2:16:ASN:HD21	1:6:1095:U:H1'	369.54	0.53
1:6:1738:U:H2'	1:6:1739:C:C6	2.44	0.53
12:C0:21:VAL:HB	12:C0:66:TYR:HB2	1.90	0.53
13:C1:17:PRO:HG3	13:C1:63:LEU:HD11	2.31	0.53
20:C8:101:LEU:O	20:C8:104:ASN:HB3	2.09	0.53
27:D5:61:SER:H	27:D5:64:VAL:HB	1.73	0.53
40:L3:108:GLU:HG2	40:L3:109:HIS:CE1	3.15	0.53
40:L3:37:ARG:HA	40:L3:186:GLY:HA2	3.38	0.53
41:L4:265:GLU:CD	41:L4:265:GLU:H	2.11	0.53
41:L4:289:ILE:O	41:L4:292:SER:OG	3.66	0.53
42:L5:184:ASP:OD2	42:L5:187:THR:OG1	5.16	0.53
45:L8:108:ARG:NE	45:L8:112:GLU:OE2	2.35	0.53
47:M0:90:ARG:HG3	47:M0:137:SER:HB2	1.89	0.53
49:M3:32:LYS:HA	49:M3:35:ARG:NH1	2.57	0.53
58:N2:104:ARG:NH1	58:N2:106:ALA:HB2	2.89	0.53
66:O0:100:ILE:HG13	66:O0:101:LEU:HD22	5.21	0.53
4:S2:218:ILE:O	4:S2:221:THR:OG1	3.86	0.53
36:1:13:A:H4'	61:N5:39:LYS:HG3	1.90	0.53
36:1:887:G:H2'	36:1:888:A:C8	2.43	0.53
36:5:1464:G:N2	36:5:1466:G:H3'	2.23	0.53
36:5:1563:C:H2'	36:5:1564:U:O4'	2.08	0.53
36:5:2261:G:H21	36:5:2262:A:N6	2.05	0.53
36:5:1899:G:N1	36:5:2335:G:OP2	2.39	0.53
36:5:255:A:H2'	36:5:256:G:H8	1.74	0.53
36:5:3065:G:H2'	36:5:3066:U:O4'	2.08	0.53
1:6:879:G:H2'	1:6:880:C:C6	2.43	0.53
18:C6:38:LEU:O	18:C6:40:GLU:N	2.68	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:D2:6:VAL:HG13	24:D2:29:PRO:HD2	2.13	0.53
26:D4:38:ASP:HA	26:D4:41:ARG:HD3	1.89	0.53
41:L4:283:THR:HG22	41:L4:285:ASP:N	2.23	0.53
49:M3:157:ARG:HD3	64:N8:101:VAL:HG21	5.28	0.53
55:M9:115:ILE:HG13	55:M9:119:LEU:HD23	1.90	0.53
56:N0:79:VAL:HG21	56:N0:106:LEU:HD11	2.90	0.53
70:O4:80:ARG:HB2	70:O4:85:VAL:HG13	5.45	0.53
76:Q0:81:SER:OG	76:Q0:82:LEU:N	2.40	0.53
2:S0:13:ASP:CG	2:S0:179:ARG:HH22	4.69	0.53
3:S1:60:ALA:O	3:S1:64:ARG:NE	7.34	0.53
3:S1:97:LEU:HD22	3:S1:232:HIS:ND1	5.44	0.53
5:S3:104:SER:OG	5:S3:105:MET:N	2.41	0.53
6:S4:49:ARG:HH11	6:S4:50:ASN:HD21	1.55	0.53
7:S5:32:GLU:HG3	7:S5:33:VAL:HG23	4.68	0.53
10:S8:184:LEU:HB3	10:S8:189:LEU:HB2	1.96	0.53
36:1:2179:C:O3'	39:L2:174:ARG:NH2	2.41	0.53
36:1:3015:G:N2	36:1:3040:A:H1'	2.24	0.53
1:2:610:G:H2'	1:2:614:C:C6	2.43	0.53
44:L7:196:LYS:HE2	36:5:1100:U:OP2	244.54	0.53
36:5:1817:G:O2'	36:5:1818:U:OP2	2.20	0.53
36:5:629:U:H2'	36:5:630:A:C8	2.42	0.53
1:6:230:C:N3	1:6:235:G:N2	2.44	0.53
19:C7:57:LEU:O	19:C7:61:ILE:N	2.92	0.53
22:D0:23:ARG:HD2	22:D0:90:TYR:CD1	2.42	0.53
23:D1:41:GLU:H	23:D1:41:GLU:CD	2.11	0.53
26:D4:125:LEU:HG	26:D4:128:LYS:HD3	6.63	0.53
39:L2:30:ARG:NH2	39:L2:33:ASP:OD2	2.46	0.53
40:L3:19:ARG:HB3	40:L3:232:ARG:HH12	1.96	0.53
40:L3:32:PHE:CD1	40:L3:182:GLN:HB3	2.43	0.53
44:L7:121:LYS:HD3	44:L7:121:LYS:O	4.51	0.53
44:L7:92:ILE:HD11	54:M8:4:ASP:HB2	1.90	0.53
47:M0:21:ARG:O	47:M0:24:ARG:NH1	2.34	0.53
50:M4:21:VAL:HA	50:M4:66:THR:HG23	2.44	0.53
57:N1:69:LYS:HZ2	57:N1:70:SER:HB3	5.47	0.53
59:N3:54:LEU:HD11	59:N3:119:GLY:HA3	1.90	0.53
67:O1:77:ARG:HD2	67:O1:89:LEU:HD23	1.90	0.53
69:O3:20:LYS:HG2	69:O3:21:ARG:HG3	1.89	0.53
72:O6:33:ALA:HB1	72:O6:38:LYS:HD2	4.59	0.53
10:S8:8:ARG:NE	10:S8:21:PHE:HD1	2.05	0.53
11:S9:93:LEU:O	11:S9:96:VAL:HG22	2.07	0.53
1:2:1400:A:H4'	19:C7:60:ARG:HH22	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:14:C:OP1	4:S2:164:SER:OG	2.18	0.53
1:2:46:A:N1	1:2:432:G:O2'	2.37	0.53
38:4:53:A:H5'	38:4:54:A:OP2	2.07	0.53
36:5:1483:G:C8	36:5:1485:G:C8	2.96	0.53
36:5:3242:G:H5'	36:5:3245:A:C8	2.39	0.53
36:5:3251:U:H2'	36:5:3252:G:C8	2.44	0.53
36:5:955:U:H2'	36:5:956:U:C6	2.43	0.53
1:6:1042:G:H22	1:6:1076:A:H2	1.57	0.53
1:6:1564:U:H2'	1:6:1565:C:C6	2.44	0.53
1:6:1591:C:H2'	1:6:1592:A:C8	2.44	0.53
1:6:20:G:H5'	1:6:571:G:C8	2.42	0.53
1:6:647:G:H22	1:6:687:G:N2	2.06	0.53
36:5:409:A:H61	38:8:15:G:H1'	1.73	0.53
12:C0:87:UNK:O	12:C0:89:UNK:N	4.37	0.53
18:C6:29:ILE:HA	18:C6:65:ILE:HB	2.82	0.53
18:C6:16:ALA:HB2	18:C6:72:GLY:HA3	1.89	0.53
19:C7:29:GLN:HA	19:C7:32:LYS:HE2	2.73	0.53
27:D5:91:PRO:HB3	27:D5:101:TYR:CE1	3.44	0.53
33:E1:147:VAL:HG23	33:E1:148:TYR:CG	2.42	0.53
39:L2:83:HIS:NE2	39:L2:86:GLN:HB2	2.24	0.53
42:L5:279:LYS:NZ	42:L5:282:ARG:HH12	2.97	0.53
42:L5:60:ILE:H	42:L5:80:SER:HB2	1.73	0.53
43:L6:62:THR:OG1	43:L6:78:ARG:HD3	2.08	0.53
43:L6:7:PRO:HD3	68:O2:74:PHE:HE1	1.99	0.53
50:M4:24:LYS:NZ	50:M4:61:GLY:O	2.24	0.53
52:M6:148:LYS:O	52:M6:150:GLU:N	2.34	0.53
55:M9:104:ARG:NH1	36:5:1949:G:H5''	215.40	0.53
56:N0:28:ARG:HH11	56:N0:99:ARG:NE	2.46	0.53
3:S1:35:PRO:HB3	3:S1:231:LEU:HD21	5.60	0.53
8:S6:136:LYS:HG3	8:S6:173:PRO:HB3	2.60	0.53
10:S8:58:LEU:O	10:S8:59:ARG:HB2	2.07	0.53
11:S9:6:ARG:HB2	11:S9:6:ARG:HH11	3.09	0.53
36:1:2228:A:H2'	36:1:2229:A:H8	1.72	0.53
36:1:3006:A:H2'	36:1:3007:U:O4'	2.07	0.53
36:1:712:G:N2	36:1:754:G:O3'	2.41	0.53
1:2:1046:G:OP1	3:S1:157:GLN:NE2	2.27	0.53
37:3:97:A:H2'	37:3:98:C:C6	2.43	0.53
38:4:83:C:H1'	38:4:85:G:H21	1.73	0.53
1:6:1068:C:H2'	1:6:1069:A:C8	2.43	0.53
21:C9:122:ARG:NH2	1:6:1500:C:OP1	419.05	0.53
20:C8:132:ARG:CZ	1:6:1544:U:H4'	344.65	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:683:C:H3'	1:6:684:A:H5''	1.91	0.53
1:6:870:C:H2'	1:6:871:G:H8	1.73	0.53
38:8:155:A:C2	38:8:156:U:H1'	2.43	0.53
17:C5:87:PRO:HD3	17:C5:112:LEU:HD22	1.90	0.53
25:D3:69:ARG:NH1	25:D3:116:ASP:OD1	2.42	0.53
26:D4:92:VAL:HG21	26:D4:99:LYS:HE3	1.90	0.53
39:L2:219:ILE:HG21	39:L2:223:SER:HB3	2.24	0.53
40:L3:139:GLN:C	40:L3:141:GLY:H	2.58	0.53
41:L4:217:LYS:HA	41:L4:220:ARG:HG2	4.47	0.53
47:M0:81:GLY:O	47:M0:83:ASP:N	2.96	0.53
53:M7:48:LEU:HB3	53:M7:88:VAL:HG13	1.91	0.53
53:M7:66:SER:HB2	36:5:1448:U:H5''	172.39	0.53
72:O6:43:LEU:O	72:O6:47:ILE:HG13	2.09	0.53
7:S5:225:ARG:HH22	30:D8:57:MET:HB2	3.74	0.53
9:S7:12:ALA:HB3	9:S7:13:PRO:HD3	1.92	0.53
9:S7:172:VAL:HG12	9:S7:176:LEU:HD12	2.82	0.53
36:1:821:U:H2'	36:1:822:G:C8	2.43	0.53
1:2:1191:U:H5'	18:C6:143:ARG:HH11	1.73	0.53
1:2:1394:G:H1	1:2:1404:C:H42	1.56	0.53
1:2:278:U:OP1	1:2:279:G:N2	2.42	0.53
38:4:34:U:O2'	38:4:35:C:OP2	2.26	0.53
36:5:2180:G:H2'	36:5:2181:C:C6	2.44	0.53
70:O4:96:GLU:OE1	36:5:2555:G:N1	213.56	0.53
1:6:514:G:O2'	1:6:515:A:H8	1.89	0.53
26:D4:11:LYS:NZ	1:6:775:G:O6	413.97	0.53
71:O5:56:THR:HG23	38:8:60:U:H3	50.88	0.53
13:C1:3:THR:HG22	13:C1:4:GLU:H	2.63	0.53
21:C9:108:LEU:HA	21:C9:111:ILE:HG22	1.90	0.53
54:M8:19:PRO:HD3	54:M8:30:VAL:HG21	2.13	0.53
64:N8:85:ASP:OD2	64:N8:86:LYS:N	3.78	0.53
4:S2:118:ALA:HB3	4:S2:124:ALA:HB2	2.02	0.53
5:S3:21:LEU:HD22	5:S3:25:PHE:CE2	2.44	0.53
5:S3:79:TYR:CD2	5:S3:84:ILE:HG13	2.44	0.53
6:S4:205:PHE:HB3	6:S4:221:ARG:HD2	1.90	0.53
8:S6:137:ARG:NH2	8:S6:177:ARG:HE	2.07	0.53
11:S9:112:GLN:HA	11:S9:112:GLN:HE21	1.73	0.53
11:S9:96:VAL:HA	11:S9:99:LEU:HB2	2.96	0.53
36:1:1783:U:H2'	36:1:1784:G:C8	2.44	0.53
36:1:3393:U:H2'	36:1:3394:U:C6	2.43	0.53
36:1:3:U:H2'	36:1:4:U:C6	2.44	0.53
1:2:1101:G:O2'	24:D2:4:SER:OG	2.14	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1523:G:N7	21:C9:68:ARG:NH1	2.56	0.53
38:4:91:C:H2'	38:4:92:A:H8	1.73	0.53
36:5:109:A:H4'	36:5:110:G:OP1	2.08	0.53
36:5:982:C:H42	36:5:1101:G:H1	1.57	0.53
36:5:2103:U:H2'	36:5:2104:A:C8	2.44	0.53
36:5:2695:A:HO2'	36:5:2696:A:H8	1.56	0.53
36:5:3163:A:N6	36:5:3288:G:O6	2.41	0.53
1:6:1429:G:H2'	1:6:1430:U:C6	2.44	0.53
1:6:1535:U:H4'	1:6:1535:U:OP1	2.08	0.53
21:C9:73:VAL:HG22	21:C9:105:LEU:HD12	1.90	0.53
6:S4:92:LEU:HG	26:D4:17:LEU:HD21	1.90	0.53
40:L3:144:ILE:HG22	40:L3:148:LEU:HD22	3.33	0.53
40:L3:289:ASP:OD2	40:L3:290:ASP:N	5.61	0.53
42:L5:82:GLU:OE2	42:L5:108:ARG:NH1	3.84	0.53
45:L8:100:GLU:OE2	45:L8:108:ARG:NH1	2.42	0.53
45:L8:61:GLN:HA	45:L8:64:ILE:HD12	1.91	0.53
55:M9:81:ARG:HG2	55:M9:88:ARG:HH21	4.24	0.53
68:O2:32:TRP:CZ2	68:O2:53:PRO:HD2	2.51	0.53
36:1:2223:A:H4'	72:O6:80:PHE:HE2	1.74	0.53
2:S0:13:ASP:HA	2:S0:16:LEU:HD12	1.91	0.53
2:S0:38:PHE:HB2	2:S0:49:ASN:HB2	1.91	0.53
8:S6:186:ARG:O	8:S6:190:GLN:HG2	2.09	0.53
36:1:1522:U:OP2	61:N5:121:LYS:NZ	2.39	0.53
36:1:1661:G:H2'	36:1:1662:G:C8	2.44	0.53
36:1:2771:U:H2'	36:1:2772:C:C2	2.44	0.53
36:1:3023:U:H2'	36:1:3024:A:C8	2.42	0.53
1:2:1402:G:P	19:C7:10:LYS:HZ1	2.32	0.53
1:2:704:C:O2	1:2:705:U:O2'	2.26	0.53
37:3:106:U:H2'	37:3:107:C:C6	2.44	0.53
78:Q2:63:LYS:NZ	36:5:2761:G:N7	212.13	0.53
76:Q0:125:LYS:NZ	36:5:2898:G:N7	327.08	0.53
64:N8:8:THR:HG21	36:5:662:U:OP1	149.75	0.53
1:6:1003:A:H4'	1:6:1004:U:O5'	2.08	0.53
25:D3:7:ARG:HD2	1:6:1102:G:OP2	351.38	0.53
24:D2:37:PHE:CZ	24:D2:103:ILE:HD12	3.62	0.53
42:L5:78:ALA:HB3	42:L5:105:ILE:HB	2.92	0.53
44:L7:103:LEU:HG	44:L7:130:ILE:HG12	4.83	0.53
46:L9:163:GLN:HB3	46:L9:166:ARG:HH11	1.73	0.53
49:M3:16:LYS:O	49:M3:18:TRP:N	3.00	0.53
52:M6:125:ARG:HG3	52:M6:129:LEU:HD22	2.65	0.53
55:M9:6:THR:HG22	55:M9:10:LEU:HD22	3.12	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:N0:155:ARG:HH21	56:N0:172:TYR:H	5.87	0.53
59:N3:39:VAL:O	59:N3:42:SER:OG	3.23	0.53
61:N5:115:ARG:HD2	61:N5:121:LYS:HB2	1.91	0.53
78:Q2:28:TYR:HB3	78:Q2:69:VAL:HB	1.89	0.53
34:SR:52:GLN:HG2	34:SR:53:LYS:HG2	2.85	0.53
36:1:1397:C:H2'	36:1:1398:U:C6	2.44	0.53
36:1:590:G:C2	36:1:610:G:H2'	2.43	0.53
1:2:442:C:O2'	1:2:525:A:N1	2.41	0.53
1:2:683:C:H2'	1:2:684:A:C8	2.44	0.53
1:2:714:G:N2	1:2:725:U:O4	2.40	0.53
36:5:651:G:C6	36:5:652:G:C6	2.97	0.53
1:6:1237:G:H2'	1:6:1238:A:H8	1.72	0.53
31:D9:14:TYR:HH	1:6:1553:G:HO2'	403.76	0.53
8:S6:87:ARG:NH2	1:6:161:U:OP2	315.10	0.53
14:C2:103:LEU:HD23	14:C2:115:VAL:HA	2.50	0.53
20:C8:102:ALA:O	20:C8:105:VAL:HG12	2.08	0.53
33:E1:105:TYR:HB3	33:E1:117:LEU:HB2	1.91	0.53
39:L2:3:ARG:HD3	36:5:911:C:N4	179.35	0.53
49:M3:57:VAL:HG13	49:M3:147:ILE:HD13	1.91	0.53
51:M5:58:GLY:HA3	51:M5:142:ILE:HD11	1.90	0.53
52:M6:177:LYS:O	52:M6:181:ALA:N	2.40	0.53
70:O4:41:ARG:HG2	70:O4:56:THR:HG21	1.90	0.53
75:O9:28:ARG:NH1	75:O9:36:ARG:HH11	6.49	0.53
79:Q3:7:LYS:O	79:Q3:27:LYS:NZ	2.41	0.53
3:S1:34:ALA:HB3	3:S1:41:ARG:HA	1.89	0.53
7:S5:89:ILE:HD12	7:S5:90:ILE:HG12	1.91	0.53
11:S9:54:ARG:HA	11:S9:57:ARG:HB3	5.00	0.53
36:1:2544:U:H2'	36:1:2545:C:C6	2.44	0.53
1:2:1186:U:OP1	1:2:1456:C:O2'	2.23	0.53
1:2:1520:U:OP2	21:C9:75:LYS:NZ	2.36	0.53
65:N9:50:THR:HG23	36:5:1073:U:H1'	204.57	0.53
21:C9:89:ARG:NH2	1:6:1562:G:OP1	375.13	0.53
12:C0:10:LYS:HD3	12:C0:36:ASP:HB3	1.91	0.53
13:C1:8:GLN:OE1	13:C1:14:GLN:N	2.33	0.53
14:C2:66:VAL:HG11	14:C2:71:ILE:HD12	1.91	0.53
17:C5:87:PRO:HA	17:C5:90:ILE:HG13	2.21	0.53
21:C9:28:LEU:HD13	21:C9:30:VAL:HG22	1.90	0.53
22:D0:50:LEU:HD23	22:D0:93:LEU:HD22	1.91	0.53
39:L2:116:VAL:HA	39:L2:164:GLY:HA2	2.30	0.53
39:L2:208:ASP:OD2	36:5:912:G:N1	186.72	0.53
39:L2:65:ASP:HB3	39:L2:68:LYS:O	2.50	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:L3:112:ASP:O	40:L3:116:ARG:HB2	2.09	0.53
40:L3:274:SER:OG	36:5:3139:A:OP1	228.05	0.53
41:L4:181:VAL:HG11	41:L4:224:GLY:HA3	2.84	0.53
47:M0:55:ASN:HA	47:M0:131:ILE:HG23	2.24	0.53
53:M7:36:ILE:HG12	53:M7:44:ALA:HB1	1.91	0.53
68:O2:81:ASP:O	68:O2:84:THR:OG1	2.88	0.53
76:Q0:125:LYS:HG3	36:5:2897:A:H5''	325.92	0.53
77:Q1:2:ARG:HG3	77:Q1:4:LYS:H	2.01	0.53
2:S0:111:ILE:HG21	1:6:1293:U:H1'	418.24	0.53
7:S5:184:PHE:CD2	7:S5:185:ARG:HG2	2.44	0.53
8:S6:105:ASP:N	8:S6:105:ASP:OD2	2.40	0.53
1:2:1681:A:H8	8:S6:66:GLY:HA3	1.74	0.53
36:1:1055:A:N3	37:3:81:U:O2'	2.40	0.52
36:1:1169:A:H5''	44:L7:219:LYS:HD2	1.91	0.52
36:1:1612:A:H5''	74:O8:51:LEU:HD12	1.90	0.52
36:1:2407:C:O2	36:1:2818:U:N3	2.33	0.52
36:1:837:A:OP1	79:Q3:5:THR:OG1	2.27	0.52
36:5:1232:C:C5	36:5:1261:G:H2'	2.44	0.52
36:5:2546:C:H2'	36:5:2547:A:H8	1.75	0.52
36:5:3156:U:O2'	36:5:3157:U:O4'	2.27	0.52
68:O2:26:HIS:HB2	36:5:655:C:H5''	160.97	0.52
1:6:1081:A:H1'	1:6:1082:C:H5	1.73	0.52
13:C1:57:LYS:NZ	1:6:326:G:OP1	289.91	0.52
12:C0:31:LYS:H	12:C0:38:LYS:HA	4.48	0.52
12:C0:77:ARG:HD3	12:C0:84:UNK:HA	1.91	0.52
13:C1:74:THR:HG22	13:C1:122:ILE:HG12	4.28	0.52
25:D3:19:ARG:NH1	1:6:610:G:H21	341.74	0.52
40:L3:220:VAL:O	40:L3:334:ARG:NH1	2.42	0.52
41:L4:156:LEU:HD23	41:L4:159:ILE:HD12	1.90	0.52
43:L6:105:TYR:HE1	43:L6:134:ARG:HD2	1.74	0.52
47:M0:77:THR:HG22	47:M0:82:ARG:HA	2.11	0.52
48:M1:16:LYS:HG2	48:M1:130:VAL:HG13	2.99	0.52
46:L9:47:LYS:HZ2	50:M4:6:ILE:H	1.57	0.52
52:M6:89:SER:O	52:M6:92:THR:OG1	2.21	0.52
54:M8:63:SER:OG	54:M8:64:VAL:N	3.03	0.52
54:M8:96:PHE:CD2	54:M8:97:PRO:HD2	2.44	0.52
72:O6:78:GLY:O	72:O6:79:SER:HB3	3.81	0.52
74:O8:11:PHE:HE1	74:O8:65:LEU:HD21	2.76	0.52
3:S1:81:PHE:HD2	3:S1:82:ARG:HG3	1.74	0.52
4:S2:53:ILE:HD11	4:S2:73:LEU:HB2	2.56	0.52
5:S3:210:GLU:OE2	19:C7:19:ARG:NH1	3.78	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:133:THR:HG21	9:S7:159:VAL:HG13	4.99	0.52
10:S8:76:THR:HG21	10:S8:104:ILE:HB	3.97	0.52
35:SM:84:LYS:HD3	35:SM:84:LYS:H	1.74	0.52
36:1:2534:G:H22	36:1:2545:C:H42	1.58	0.52
36:1:2768:U:H2'	36:1:2769:A:C8	2.43	0.52
36:1:3116:G:N2	36:1:3116:G:OP1	2.32	0.52
36:1:3151:U:OP1	40:L3:128:LYS:NZ	2.36	0.52
36:1:300:G:H1	36:1:315:C:H42	1.57	0.52
36:1:501:A:H2'	36:1:502:U:C6	2.45	0.52
36:1:874:U:N3	36:1:2978:U:OP1	2.39	0.52
36:1:976:U:P	54:M8:144:ARG:HH22	2.31	0.52
1:2:1339:C:O2'	1:2:1341:A:N7	2.27	0.52
50:M4:42:LYS:HD3	36:5:1185:C:OP1	305.56	0.52
36:5:1222:G:H1'	36:5:1285:G:H22	1.73	0.52
45:L8:32:LYS:HB2	36:5:2561:A:N1	203.31	0.52
1:6:1533:C:H4'	1:6:1539:G:N1	2.24	0.52
1:2:926:A:H2	16:C4:125:SER:HB3	1.72	0.52
19:C7:65:PRO:HB3	19:C7:74:GLN:NE2	3.48	0.52
20:C8:91:ASP:C	20:C8:93:THR:H	2.12	0.52
23:D1:71:ARG:HG2	23:D1:83:TRP:CH2	2.45	0.52
39:L2:177:LYS:HD2	79:Q3:69:TYR:CE1	3.16	0.52
40:L3:22:ALA:H	40:L3:272:TYR:HD1	1.55	0.52
40:L3:44:THR:HA	40:L3:340:LYS:HD3	5.03	0.52
40:L3:66:LYS:HB3	40:L3:66:LYS:NZ	2.25	0.52
49:M3:75:PHE:O	49:M3:79:GLU:HB2	2.08	0.52
52:M6:10:ASP:OD2	52:M6:37:ARG:NH2	2.98	0.52
57:N1:14:MET:CE	57:N1:55:LYS:HB2	2.64	0.52
77:Q1:2:ARG:HB3	77:Q1:5:TRP:CD1	2.43	0.52
2:S0:72:ASP:OD2	2:S0:72:ASP:N	3.01	0.52
8:S6:31:ARG:HD2	8:S6:34:GLN:HE22	1.74	0.52
9:S7:46:ILE:HD11	9:S7:58:LEU:HB3	1.90	0.52
1:2:396:G:O6	10:S8:26:LYS:NZ	2.39	0.52
36:1:2157:G:N7	39:L2:152:SER:OG	2.42	0.52
36:1:3072:C:H2'	36:1:3073:A:O4'	2.08	0.52
36:5:1025:A:H3'	36:5:1026:A:H4'	1.90	0.52
70:O4:10:ARG:O	36:5:1488:G:O2'	139.28	0.52
39:L2:241:ARG:HH22	36:5:2156:C:P	213.79	0.52
36:5:3343:G:H21	36:5:3362:A:H2	1.55	0.52
4:S2:88:LYS:HE3	1:6:1301:U:H5'	382.93	0.52
1:6:1350:U:H2'	1:6:1351:G:H8	1.75	0.52
1:6:292:U:H2'	1:6:293:U:C6	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:C7:29:GLN:OE1	19:C7:32:LYS:NZ	3.59	0.52
24:D2:119:LYS:HB3	24:D2:121:VAL:HG23	1.89	0.52
28:D6:44:ILE:HD13	28:D6:65:PRO:HB2	6.86	0.52
39:L2:174:ARG:HA	79:Q3:69:TYR:CE2	2.79	0.52
39:L2:49:VAL:HG22	39:L2:50:HIS:H	1.74	0.52
41:L4:141:ARG:CZ	41:L4:180:LYS:HG3	3.49	0.52
47:M0:43:VAL:HA	47:M0:139:ARG:HH22	1.74	0.52
49:M3:107:GLU:OE1	72:O6:18:THR:OG1	3.30	0.52
53:M7:127:ARG:O	53:M7:139:TYR:N	3.15	0.52
57:N1:56:PHE:CZ	57:N1:78:LYS:HD3	2.86	0.52
59:N3:90:GLY:O	60:N4:16:GLY:HA2	2.38	0.52
36:1:718:G:OP1	64:N8:117:ARG:NH2	2.42	0.52
74:O8:11:PHE:HD1	74:O8:12:LEU:HD22	1.74	0.52
3:S1:77:GLU:O	3:S1:80:SER:OG	4.36	0.52
5:S3:64:ARG:HA	5:S3:67:ASN:HB2	3.27	0.52
9:S7:154:LEU:HD21	9:S7:183:PHE:CD1	2.41	0.52
10:S8:26:LYS:HE3	1:6:396:G:O6	303.64	0.52
34:SR:88:THR:HG22	34:SR:104:VAL:HG22	1.91	0.52
34:SR:203:THR:OG1	34:SR:204:ALA:N	2.69	0.52
36:1:950:G:N1	36:1:1368:U:OP2	2.31	0.52
36:1:153:U:HO2'	36:1:158:G:HO2'	1.53	0.52
36:1:1940:G:H2'	36:1:1941:C:O4'	2.10	0.52
36:1:42:C:O2'	36:1:43:A:H5'	2.08	0.52
1:2:206:A:H1'	1:2:262:U:C2	2.45	0.52
1:2:696:C:O3'	1:2:697:C:H2'	2.09	0.52
42:L5:140:ARG:HD3	36:5:1080:A:OP1	227.38	0.52
36:5:1773:C:H2'	36:5:1774:C:H6	1.73	0.52
36:5:2709:C:H2'	36:5:2710:C:H6	1.74	0.52
1:6:1767:G:OP1	1:6:1770:U:H4'	2.10	0.52
14:C2:124:LYS:O	14:C2:126:TRP:N	2.36	0.52
15:C3:63:ALA:O	15:C3:67:THR:HG23	4.30	0.52
23:D1:8:LEU:HD13	23:D1:9:VAL:H	2.89	0.52
28:D6:87:ARG:HD3	1:6:1796:C:OP1	346.01	0.52
29:D7:36:LYS:HB3	29:D7:43:ILE:HG22	1.91	0.52
30:D8:11:LYS:HZ3	30:D8:33:LEU:HD21	1.74	0.52
39:L2:201:GLY:HA2	39:L2:204:MET:SD	2.54	0.52
41:L4:206:LEU:HD23	41:L4:226:GLU:HB2	1.91	0.52
41:L4:295:ILE:O	41:L4:299:ILE:HG12	2.36	0.52
42:L5:210:GLU:O	42:L5:214:ASP:HB2	2.95	0.52
44:L7:140:SER:N	44:L7:237:ASN:OD1	2.32	0.52
46:L9:166:ARG:NH2	46:L9:168:ARG:HH12	13.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
66:O0:99:ASP:OD1	66:O0:103:THR:OG1	4.67	0.52
67:O1:44:MET:HB2	67:O1:46:THR:HG22	1.90	0.52
70:O4:21:LYS:HG2	70:O4:33:GLN:HB3	1.91	0.52
36:1:1494:U:OP1	75:O9:44:TRP:HB3	2.10	0.52
2:S0:172:LEU:HD22	2:S0:176:LEU:HG	2.37	0.52
3:S1:141:ALA:HB1	3:S1:207:LEU:HD22	3.76	0.52
3:S1:62:LYS:O	3:S1:64:ARG:N	2.39	0.52
5:S3:132:LYS:HG3	5:S3:156:PHE:HB3	2.02	0.52
10:S8:44:HIS:O	10:S8:56:ARG:N	2.88	0.52
36:1:1278:A:O2'	36:1:1279:C:O5'	2.27	0.52
36:1:2376:G:H2'	36:1:2377:G:C8	2.44	0.52
36:1:2675:C:N4	48:M1:22:SER:HB3	2.24	0.52
36:1:2916:U:H2'	36:1:2917:G:H8	1.75	0.52
36:1:28:C:O2'	36:1:61:A:N3	2.37	0.52
1:2:1089:U:O2'	1:2:1093:A:N1	2.36	0.52
1:2:513:U:H2'	1:2:514:G:C8	2.45	0.52
37:3:19:C:H42	37:3:60:G:H1	1.57	0.52
39:L2:70:ARG:NH1	36:5:1580:A:OP1	174.29	0.52
36:5:2256:A:OP2	36:5:2256:A:H2'	2.09	0.52
36:5:2716:U:C2'	36:5:2717:U:H5'	2.40	0.52
52:M6:161:LYS:HD3	36:5:3182:G:H4'	287.11	0.52
1:6:1776:A:H2'	1:6:1777:G:C8	2.45	0.52
4:S2:197:TYR:HD1	1:6:2:A:H2'	392.92	0.52
20:C8:120:ARG:HE	35:SM:61:ILE:HG21	5.36	0.52
1:2:778:G:H1	26:D4:10:ARG:HG2	1.74	0.52
39:L2:89:TYR:N	39:L2:100:ASN:OD1	2.97	0.52
39:L2:243:THR:OG1	36:5:2244:A:H5"	227.77	0.52
39:L2:3:ARG:HD3	36:5:911:C:H42	179.72	0.52
45:L8:51:LYS:HG3	36:5:2523:A:C5	165.10	0.52
46:L9:8:GLN:HG2	46:L9:68:LEU:HD13	2.68	0.52
51:M5:27:VAL:HB	51:M5:122:ASN:ND2	2.24	0.52
51:M5:183:THR:HB	51:M5:187:ARG:HB2	1.91	0.52
61:N5:45:LYS:HD3	71:O5:75:TYR:CE2	3.26	0.52
63:N7:133:LYS:HE3	63:N7:135:ARG:HD3	1.90	0.52
66:O0:43:ILE:HD11	66:O0:92:ILE:HD11	1.91	0.52
70:O4:74:ARG:NH2	70:O4:82:ALA:HB2	2.25	0.52
64:N8:149:ALA:HB3	72:O6:15:LYS:HD2	1.91	0.52
4:S2:238:SER:C	4:S2:240:LEU:H	2.11	0.52
6:S4:176:ASP:OD2	6:S4:176:ASP:N	3.04	0.52
6:S4:191:ARG:CZ	6:S4:245:LYS:HD3	2.99	0.52
6:S4:37:LYS:O	6:S4:40:GLU:N	3.22	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:151:LYS:HD2	9:S7:182:VAL:HG11	1.92	0.52
34:SR:236:ALA:O	34:SR:261:LYS:NZ	4.14	0.52
36:1:118:U:H3	36:1:122:A:H5'	1.75	0.52
36:1:1222:G:O2'	36:1:1285:G:N1	2.43	0.52
36:1:1744:G:H2'	36:1:1745:C:C6	2.44	0.52
36:1:238:A:C2'	36:1:239:G:H5'	2.40	0.52
36:1:2969:A:N7	39:L2:215:ASN:ND2	2.57	0.52
36:1:3157:U:H4'	36:1:3158:G:H5'	1.91	0.52
1:2:1104:U:H5	25:D3:4:GLY:HA2	1.75	0.52
1:2:354:C:H5''	10:S8:16:ALA:HB2	1.90	0.52
1:2:639:U:OP1	9:S7:117:THR:OG1	2.18	0.52
1:2:70:C:H2'	1:2:71:A:O4'	2.10	0.52
1:2:716:C:H41	1:2:723:G:H21	1.56	0.52
37:3:60:G:H2'	37:3:61:G:C8	2.45	0.52
38:4:71:A:N1	38:4:82:U:O2'	2.37	0.52
36:5:1564:U:H2'	36:5:1565:G:C8	2.44	0.52
55:M9:9:ARG:NH2	36:5:1603:A:H5'	110.30	0.52
1:6:138:A:H5''	1:6:138:A:N3	2.24	0.52
8:S6:66:GLY:HA3	1:6:1681:A:H1'	273.91	0.52
1:6:913:G:H3'	1:6:914:G:H5'	1.92	0.52
45:L8:63:LYS:NZ	38:8:152:G:O2'	147.94	0.52
12:C0:44:LYS:NZ	12:C0:47:GLN:OE1	2.30	0.52
13:C1:121:ASP:OD1	13:C1:145:ALA:HB2	6.13	0.52
20:C8:131:LEU:HA	20:C8:145:ARG:HH12	1.74	0.52
21:C9:86:ARG:HB2	21:C9:89:ARG:HB2	1.90	0.52
23:D1:32:VAL:HG12	23:D1:55:LEU:HB2	2.07	0.52
24:D2:99:PHE:H	24:D2:99:PHE:HD1	2.59	0.52
39:L2:174:ARG:NH2	36:5:2179:C:O3'	213.51	0.52
44:L7:123:THR:HA	44:L7:126:LEU:HB2	1.91	0.52
46:L9:111:PHE:HD1	46:L9:127:PRO:HA	2.14	0.52
46:L9:137:SER:HB2	46:L9:143:GLU:HB3	2.71	0.52
47:M0:80:SER:HB3	47:M0:147:VAL:HG11	1.92	0.52
48:M1:117:ASP:OD2	48:M1:119:SER:OG	2.18	0.52
49:M3:61:PRO:O	49:M3:62:THR:OG1	2.40	0.52
42:L5:34:LYS:HD2	57:N1:30:TYR:CZ	2.45	0.52
57:N1:78:LYS:NZ	36:5:2728:G:O6	219.13	0.52
64:N8:126:LYS:HB3	64:N8:148:ILE:HG21	1.92	0.52
68:O2:24:ARG:HG2	68:O2:25:TYR:CE1	3.02	0.52
3:S1:116:LYS:HZ2	3:S1:117:TRP:HZ3	1.57	0.52
5:S3:68:GLU:O	5:S3:72:LEU:N	2.94	0.52
11:S9:132:ARG:HB2	11:S9:140:ILE:HG21	4.59	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1427:U:H2'	36:1:1428:A:C8	2.44	0.52
36:1:1659:U:H2'	36:1:1660:C:C6	2.45	0.52
36:1:29:C:H4'	36:1:62:A:H4'	1.90	0.52
36:1:514:G:H21	41:L4:341:SER:HA	1.74	0.52
36:1:75:G:H5''	49:M3:58:VAL:HG13	1.92	0.52
1:2:1275:A:N3	5:S3:141:LYS:NZ	2.48	0.52
1:2:1401:A:O3'	19:C7:10:LYS:NZ	2.41	0.52
1:2:1600:A:H4'	1:2:1601:G:OP1	2.10	0.52
1:2:256:A:H2'	1:2:257:A:O4'	2.09	0.52
1:2:855:A:C2	1:2:857:U:H1'	2.44	0.52
38:4:91:C:H2'	38:4:92:A:C8	2.45	0.52
36:5:127:G:H2'	36:5:128:G:C8	2.45	0.52
75:O9:42:ARG:HH22	36:5:1494:U:P	107.99	0.52
36:5:1481:A:H2'	36:5:1858:A:N3	2.25	0.52
36:5:207:U:H2'	36:5:208:C:C6	2.45	0.52
36:5:2660:G:H1'	36:5:2744:U:H1'	1.91	0.52
36:5:305:U:O2	36:5:2784:G:N2	2.41	0.52
79:Q3:17:ARG:NH1	36:5:860:G:OP1	219.02	0.52
1:6:1016:C:H2'	1:6:1017:U:H6	1.74	0.52
1:6:209:U:H2'	1:6:210:A:C8	2.45	0.52
1:6:413:U:H2'	1:6:414:C:H6	1.75	0.52
14:C2:57:ALA:HB3	14:C2:85:LYS:HZ1	1.73	0.52
15:C3:132:VAL:HG23	15:C3:134:VAL:HG22	1.91	0.52
20:C8:134:ARG:HB3	1:6:1559:A:C8	364.48	0.52
22:D0:17:GLN:HG3	22:D0:18:GLN:HG3	8.70	0.52
27:D5:80:LEU:HD13	27:D5:101:TYR:HE2	2.65	0.52
39:L2:83:HIS:CD2	39:L2:86:GLN:HB2	2.45	0.52
42:L5:40:HIS:CD2	42:L5:42:ALA:HB3	2.44	0.52
44:L7:219:LYS:HE2	36:5:1169:A:H4'	251.58	0.52
49:M3:75:PHE:H	49:M3:97:VAL:HA	1.75	0.52
53:M7:3:ARG:HH12	36:5:398:A:N6	133.82	0.52
56:N0:133:ALA:HA	56:N0:141:LYS:HZ1	1.74	0.52
62:N6:112:ASP:HB2	62:N6:115:ARG:HB2	1.92	0.52
74:O8:63:LYS:HA	74:O8:66:ILE:HG13	3.24	0.52
5:S3:22:ASN:OD1	5:S3:34:TYR:OH	2.22	0.52
6:S4:104:ASP:HB3	6:S4:106:LYS:H	4.01	0.52
11:S9:34:PHE:CD2	11:S9:105:LEU:HD23	2.45	0.52
36:1:1355:A:H4'	36:1:1356:U:O5'	2.07	0.52
36:1:1794:G:O2'	36:1:1795:U:H5'	2.10	0.52
36:1:2338:C:H1'	59:N3:49:LEU:HD12	1.91	0.52
36:1:3114:A:O2'	46:L9:62:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:2389:C:O2'	36:1:3307:A:N1	2.41	0.52
1:2:1681:A:C8	8:S6:66:GLY:HA3	2.45	0.52
1:2:61:A:H8	1:2:269:G:HO2'	1.56	0.52
37:3:48:U:N3	37:3:49:G:N7	2.57	0.52
36:5:2524:A:H1'	36:5:2525:G:C8	2.45	0.52
36:5:999:G:C6	36:5:1000:C:N4	2.78	0.52
1:6:1435:G:H4'	1:6:1436:A:H5'	1.91	0.52
48:M1:152:HIS:HB2	37:7:56:A:H4'	326.77	0.52
37:7:92:A:C5	37:7:93:C:H1'	2.44	0.52
18:C6:55:VAL:HG21	18:C6:89:LEU:HD21	1.90	0.52
23:D1:14:PRO:HB2	23:D1:23:ILE:HG23	2.47	0.52
24:D2:86:ILE:HD12	24:D2:87:GLU:HG2	1.90	0.52
36:1:2525:G:OP2	39:L2:37:ARG:NH1	2.42	0.52
40:L3:114:VAL:O	40:L3:117:ARG:HB3	2.10	0.52
45:L8:213:LYS:O	45:L8:216:SER:OG	4.95	0.52
48:M1:39:GLN:HG3	48:M1:40:LEU:N	2.24	0.52
48:M1:8:PRO:HD2	48:M1:10:ARG:HG3	2.73	0.52
52:M6:84:LEU:HD13	52:M6:102:LEU:HD21	1.91	0.52
61:N5:63:ILE:HA	61:N5:86:VAL:HG23	2.74	0.52
63:N7:41:ALA:HB2	63:N7:77:TYR:HE1	1.73	0.52
3:S1:219:LYS:HD2	79:Q3:92:ALA:HB3	13.04	0.52
8:S6:139:ASN:HA	8:S6:142:ARG:HD2	2.55	0.52
8:S6:199:GLN:HG2	8:S6:202:ARG:HH12	1.96	0.52
9:S7:45:SER:HB3	9:S7:61:PHE:CD2	2.44	0.52
36:1:1110:U:OP1	54:M8:164:ARG:NH2	2.33	0.52
36:1:2683:U:H2'	36:1:2684:C:C6	2.45	0.52
64:N8:10:LYS:NZ	36:5:1374:G:O6	163.91	0.52
36:5:408:A:N6	38:8:15:G:H1'	2.25	0.52
36:5:638:C:H2'	36:5:639:G:H8	1.74	0.52
1:6:1325:A:H2'	1:6:1326:A:H8	1.74	0.52
10:S8:50:GLY:HA2	1:6:397:A:O3'	314.67	0.52
1:6:542:A:C8	1:6:543:C:H2'	2.44	0.52
14:C2:56:GLU:OE1	14:C2:124:LYS:NZ	3.40	0.52
20:C8:131:LEU:HA	20:C8:145:ARG:NH1	2.25	0.52
39:L2:27:ALA:HB3	39:L2:128:ARG:HH22	2.80	0.52
2:S0:11:PRO:HA	2:S0:14:ALA:HB3	3.04	0.52
2:S0:72:ASP:HB2	2:S0:118:PRO:HA	1.92	0.52
11:S9:117:GLY:O	11:S9:119:ALA:N	2.91	0.52
20:C8:146:ALA:H	35:SM:68:ARG:HH21	1.57	0.52
34:SR:123:ILE:HG22	34:SR:133:VAL:HG22	1.91	0.52
36:1:1018:G:N7	36:1:1035:G:N2	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1686:U:OP1	58:N2:42:LYS:NZ	2.27	0.52
36:1:168:U:H2'	36:1:169:U:C6	2.45	0.52
36:1:2094:C:H2'	36:1:2095:G:C8	2.45	0.52
36:1:2538:U:O2'	36:1:2541:U:O4	2.19	0.52
36:1:2883:U:H2'	36:1:2884:C:C6	2.45	0.52
36:1:3280:U:O2'	36:1:3281:U:H5''	2.10	0.52
1:2:999:U:O4	1:2:1000:C:N4	2.42	0.52
1:2:1335:U:H2'	1:2:1336:A:C8	2.44	0.52
1:2:1525:A:H2'	1:2:1526:A:C8	2.44	0.52
1:2:1602:C:H2'	1:2:1603:U:O4'	2.10	0.52
1:2:209:U:H2'	1:2:210:A:C8	2.45	0.52
1:2:248:U:H4'	13:C1:36:LYS:HD3	1.92	0.52
1:2:749:U:H5''	24:D2:83:ILE:HD12	1.91	0.52
36:5:2344:U:H2'	36:5:2345:A:C8	2.44	0.52
57:N1:54:HIS:CD2	36:5:2724:U:H4'	228.98	0.52
1:6:1032:G:H2'	1:6:1033:C:C6	2.45	0.52
1:6:1472:C:N4	1:6:1536:G:O6	2.42	0.52
1:2:1553:G:O6	17:C5:43:ARG:HD3	2.09	0.52
1:2:1312:A:N6	19:C7:2:GLY:O	2.43	0.52
23:D1:36:VAL:O	23:D1:51:VAL:N	2.83	0.52
24:D2:103:ILE:HD11	24:D2:126:LEU:HB3	3.90	0.52
24:D2:90:THR:HB	24:D2:94:LEU:HD12	2.08	0.52
26:D4:50:ALA:O	26:D4:51:GLU:HB3	2.10	0.52
27:D5:75:LEU:O	27:D5:79:ALA:N	2.92	0.52
16:C4:128:LYS:HD3	28:D6:27:SER:OG	3.33	0.52
29:D7:49:HIS:CE1	29:D7:70:LYS:HG2	2.80	0.52
44:L7:221:LYS:HB2	44:L7:227:GLY:HA3	1.92	0.52
49:M3:42:ARG:O	49:M3:46:ILE:HG12	2.30	0.52
50:M4:54:PRO:O	50:M4:56:GLN:HG2	2.10	0.52
55:M9:110:ARG:HH12	36:5:1720:U:P	231.57	0.52
55:M9:92:GLN:O	55:M9:96:ILE:HG13	2.10	0.52
56:N0:154:HIS:CE1	56:N0:170:THR:HG21	2.45	0.52
56:N0:28:ARG:HE	56:N0:99:ARG:NH2	2.49	0.52
59:N3:11:PHE:CD1	59:N3:88:ARG:HD2	3.09	0.52
64:N8:75:LEU:HB3	64:N8:118:ILE:HG23	1.92	0.52
69:O3:52:VAL:HG21	69:O3:99:ARG:NH1	2.25	0.52
72:O6:97:SER:O	72:O6:99:ARG:N	2.42	0.52
36:1:1073:U:H1'	65:N9:50:THR:HB	1.92	0.51
36:1:1915:A:H2'	36:1:1916:U:C6	2.45	0.51
36:1:3282:U:H2'	36:1:3283:U:C6	2.45	0.51
36:1:653:A:H61	36:1:1442:U:H3	1.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1413:U:H4'	1:2:1414:U:OP2	2.10	0.51
1:2:15:U:H2'	1:2:16:G:O4'	2.10	0.51
37:3:7:G:O3'	42:L5:33:ARG:NH2	2.42	0.51
36:5:1152:G:H22	36:5:1200:A:N6	2.08	0.51
36:5:1256:G:O6	36:5:1261:G:N2	2.44	0.51
11:S9:124:HIS:ND1	1:6:478:A:O2'	448.42	0.51
1:6:482:U:H2'	1:6:483:A:H8	1.75	0.51
1:6:701:U:H2'	1:6:702:G:H8	1.75	0.51
11:S9:146:PHE:HZ	1:6:765:G:H1	431.21	0.51
14:C2:136:ILE:HA	14:C2:139:HIS:HB3	1.91	0.51
15:C3:84:ILE:HD11	15:C3:89:TYR:HD2	2.07	0.51
20:C8:91:ASP:OD1	20:C8:92:ILE:N	4.46	0.51
27:D5:39:ALA:O	27:D5:72:GLY:N	2.41	0.51
33:E1:105:TYR:HA	33:E1:117:LEU:HD12	1.92	0.51
47:M0:77:THR:O	47:M0:79:VAL:N	2.34	0.51
49:M3:174:ARG:HG3	72:O6:9:ILE:HD11	5.77	0.51
56:N0:91:TYR:OH	56:N0:93:GLU:OE2	2.15	0.51
62:N6:113:LYS:HB2	38:8:84:C:H1'	19.94	0.51
78:Q2:12:CYS:SG	78:Q2:17:CYS:HB2	2.86	0.51
6:S4:98:ASN:ND2	6:S4:116:ASP:OD1	3.13	0.51
8:S6:159:ARG:HG2	8:S6:172:ALA:HB2	2.77	0.51
34:SR:129:LYS:HG2	34:SR:149:ASP:O	2.27	0.51
36:1:1488:G:H5''	36:1:1838:G:O6	2.10	0.51
36:1:1751:G:H5'	74:O8:26:LYS:NZ	2.25	0.51
36:1:2211:U:H2'	36:1:2212:C:O4'	2.10	0.51
36:1:741:U:O2'	54:M8:73:GLN:HG2	2.10	0.51
36:1:799:G:HO2'	49:M3:18:TRP:HE1	1.56	0.51
1:2:1165:G:H1	1:2:1580:C:H42	1.58	0.51
1:2:1429:G:H1'	22:D0:74:GLU:HG2	1.91	0.51
1:2:895:G:H1	1:2:917:U:H3	1.57	0.51
37:3:45:A:H2'	37:3:46:A:H8	1.73	0.51
38:4:95:G:OP1	73:O7:76:ASN:ND2	2.42	0.51
36:5:1367:G:HO2'	36:5:1368:U:H6	1.57	0.51
36:5:1879:A:OP1	36:5:1879:A:H4'	2.09	0.51
36:5:3112:G:O6	36:5:3120:C:H5''	2.10	0.51
36:5:3236:U:H2'	36:5:3237:U:H6	1.75	0.51
20:C8:139:LYS:O	1:6:1461:C:N4	343.57	0.51
1:6:1650:U:H2'	1:6:1651:A:C8	2.45	0.51
1:6:263:C:H4'	1:6:292:U:H5'	1.92	0.51
28:D6:87:ARG:HD2	1:6:1797:A:N1	345.11	0.51
36:1:2185:G:H5'	39:L2:219:ILE:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:L3:37:ARG:HH11	40:L3:191:LYS:HE3	1.75	0.51
43:L6:85:ILE:HG23	69:O3:107:ILE:HG21	2.78	0.51
49:M3:105:ASN:OD1	49:M3:107:GLU:HG2	2.98	0.51
49:M3:116:LEU:O	49:M3:120:GLN:NE2	2.42	0.51
50:M4:124:ARG:NH2	36:5:3212:C:OP2	289.10	0.51
56:N0:38:LYS:HG3	56:N0:61:ILE:HD12	3.49	0.51
60:N4:4:GLU:HG2	60:N4:30:ARG:HD2	1.92	0.51
65:N9:14:ARG:NH1	65:N9:18:ARG:HD2	2.25	0.51
2:S0:41:ARG:HB2	2:S0:47:VAL:HG23	1.91	0.51
6:S4:37:LYS:HB2	6:S4:40:GLU:HG2	1.92	0.51
36:1:1492:G:N7	75:O9:2:ALA:HB1	2.26	0.51
36:1:1744:G:H2'	36:1:1745:C:H6	1.75	0.51
36:1:2611:U:H2'	36:1:2612:U:H6	1.76	0.51
36:1:2714:G:H4'	36:1:2715:A:O5'	2.10	0.51
36:1:2947:G:C2	40:L3:250:ALA:HB1	2.45	0.51
36:1:994:G:N2	36:1:995:U:O4	2.39	0.51
1:2:788:A:H2'	6:S4:19:LEU:HD22	1.91	0.51
1:2:980:G:H4'	1:2:1776:A:H4'	1.93	0.51
36:5:1014:U:H3	36:5:1036:A:H61	1.58	0.51
36:5:1102:A:H4'	36:5:1103:A:C6	2.45	0.51
41:L4:197:ARG:NH1	36:5:1381:A:OP1	108.99	0.51
36:5:2378:C:H2'	36:5:2379:U:C6	2.45	0.51
42:L5:12:TYR:OH	36:5:2688:U:OP1	299.50	0.51
36:5:2830:G:H1	36:5:2858:U:H3	1.57	0.51
36:5:3121:U:H1'	36:5:3122:A:H5''	1.93	0.51
17:C5:77:ARG:NH1	1:6:1241:G:OP2	383.41	0.51
1:6:794:U:H4'	1:6:795:U:OP2	2.09	0.51
19:C7:104:ASN:OD1	19:C7:104:ASN:N	2.43	0.51
20:C8:131:LEU:HD23	20:C8:145:ARG:HH12	1.75	0.51
22:D0:28:SER:OG	22:D0:29:THR:N	2.44	0.51
26:D4:27:VAL:HG11	26:D4:35:VAL:HG11	1.92	0.51
40:L3:227:GLU:OE2	40:L3:270:ARG:NE	2.40	0.51
41:L4:186:LYS:HE2	36:5:1388:U:O4	117.72	0.51
46:L9:89:LYS:HB2	46:L9:183:HIS:HB3	1.91	0.51
47:M0:36:LEU:CD1	47:M0:87:LEU:HB3	3.13	0.51
48:M1:143:ARG:NH2	37:7:5:G:OP1	292.02	0.51
49:M3:14:PHE:CE1	36:5:665:A:H1'	133.75	0.51
51:M5:143:ARG:HH21	71:O5:92:LEU:HA	2.40	0.51
56:N0:24:LEU:HD21	56:N0:59:VAL:HG21	1.92	0.51
60:N4:25:ASP:OD2	60:N4:26:SER:N	4.94	0.51
63:N7:73:LYS:HG2	63:N7:74:VAL:O	4.81	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
71:O5:21:LEU:HD22	71:O5:25:LYS:HE2	4.62	0.51
38:4:41:A:O2'	73:O7:59:THR:HG22	2.09	0.51
8:S6:175:ILE:HG22	8:S6:178:LEU:HB2	3.19	0.51
36:1:1355:A:C5'	36:1:1356:U:H5	2.23	0.51
36:1:2168:A:C6	36:1:2170:U:H1'	2.45	0.51
36:1:2174:G:H8	36:1:2174:G:OP1	1.93	0.51
36:1:532:A:H2	36:1:560:G:H22	1.58	0.51
36:1:944:C:H4'	68:O2:33:ARG:NH1	2.26	0.51
1:2:1409:G:N1	1:2:1412:G:OP2	2.42	0.51
1:2:542:A:C8	1:2:543:C:H5'	2.46	0.51
36:5:123:A:C6	36:5:150:A:C5	2.98	0.51
22:D0:60:THR:HG22	1:6:1382:A:H5''	435.15	0.51
1:6:1689:A:H2'	1:6:1690:G:H8	1.74	0.51
1:6:738:G:H2'	1:6:739:G:C8	2.45	0.51
15:C3:42:ARG:C	15:C3:44:GLY:H	2.72	0.51
16:C4:132:ARG:NH1	1:6:1788:G:O5'	295.70	0.51
20:C8:116:LEU:HA	20:C8:119:ILE:HG12	5.20	0.51
24:D2:11:LEU:HD11	24:D2:37:PHE:CE2	3.20	0.51
28:D6:24:VAL:HG11	28:D6:71:LEU:HD13	1.93	0.51
40:L3:116:ARG:HB3	40:L3:175:LYS:HA	1.91	0.51
41:L4:33:ASP:OD1	41:L4:34:ILE:N	2.40	0.51
47:M0:3:ARG:CZ	47:M0:63:GLU:HG3	2.41	0.51
47:M0:72:ALA:HB2	47:M0:155:ALA:HB2	2.09	0.51
48:M1:47:GLN:HG2	48:M1:64:LYS:HE2	1.93	0.51
52:M6:15:LEU:HD21	52:M6:125:ARG:HG3	1.91	0.51
70:O4:3:GLN:NE2	70:O4:30:LEU:O	4.97	0.51
6:S4:241:GLY:O	6:S4:244:ILE:HG12	2.10	0.51
11:S9:59:LEU:HD22	11:S9:69:ARG:HA	2.80	0.51
34:SR:126:SER:OG	34:SR:127:ARG:N	2.42	0.51
34:SR:224:ASN:O	34:SR:228:LYS:HA	2.71	0.51
34:SR:31:ASN:O	34:SR:47:LEU:N	2.37	0.51
36:1:2186:U:H2'	36:1:2187:G:O4'	2.10	0.51
36:1:2707:C:H2'	36:1:2708:C:H6	1.75	0.51
36:1:3302:U:O2'	36:1:3303:G:H5'	2.10	0.51
36:1:378:A:N7	36:1:391:A:H2	2.09	0.51
1:2:1473:U:H5''	7:S5:190:ILE:HG12	1.93	0.51
1:2:194:U:HO2'	1:2:195:G:HO2'	1.56	0.51
1:2:416:A:H4'	1:2:417:A:OP2	2.10	0.51
1:2:633:U:H2'	1:2:634:G:O4'	2.11	0.51
37:3:91:G:H2'	37:3:92:A:C8	2.45	0.51
38:4:104:A:C8	38:4:105:A:C8	2.99	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:L4:220:ARG:NH1	36:5:211:A:OP1	75.13	0.51
36:5:541:U:H2'	36:5:542:G:C8	2.45	0.51
36:5:621:A:H2'	36:5:622:A:C8	2.46	0.51
1:6:1691:A:H2	1:6:1710:U:H3	1.58	0.51
1:6:205:U:H2'	1:6:206:A:C8	2.46	0.51
15:C3:34:ILE:O	15:C3:38:VAL:HG23	2.10	0.51
1:2:1550:A:P	17:C5:42:ARG:HH21	2.33	0.51
20:C8:25:ASN:HB2	27:D5:40:VAL:HB	1.92	0.51
24:D2:29:PRO:HB2	24:D2:58:SER:HB3	2.43	0.51
26:D4:92:VAL:HG21	26:D4:99:LYS:HG2	1.92	0.51
40:L3:107:ALA:HA	40:L3:199:PHE:CD2	2.46	0.51
42:L5:279:LYS:HZ2	42:L5:282:ARG:HH12	2.32	0.51
43:L6:42:LEU:HD23	43:L6:84:VAL:HG13	5.33	0.51
46:L9:87:LYS:HD2	46:L9:191:LEU:HD11	14.19	0.51
52:M6:23:VAL:HG11	52:M6:84:LEU:HD11	1.93	0.51
55:M9:127:SER:HA	55:M9:132:PHE:HD2	1.75	0.51
8:S6:160:ARG:HG3	60:N4:84:GLY:HA3	1.92	0.51
72:O6:40:VAL:O	72:O6:44:VAL:HG23	2.10	0.51
2:S0:124:THR:HB	2:S0:174:TRP:HE1	3.13	0.51
3:S1:139:ALA:HA	3:S1:212:VAL:HA	3.03	0.51
3:S1:146:GLN:O	3:S1:148:ASN:N	2.93	0.51
10:S8:30:GLY:HA2	1:6:332:U:H5'	290.33	0.51
10:S8:34:ALA:HB2	10:S8:56:ARG:HG3	1.93	0.51
34:SR:166:SER:HA	34:SR:184:ASN:HD21	1.75	0.51
36:1:2842:U:OP1	36:1:2844:C:N4	2.44	0.51
36:1:341:G:N7	41:L4:195:ARG:NH2	2.54	0.51
36:1:415:G:H2'	36:1:416:A:H8	1.74	0.51
1:2:1184:A:HO2'	1:2:1209:C:HO2'	1.55	0.51
1:2:325:G:H2'	1:2:326:G:H8	1.76	0.51
37:3:112:G:H2'	37:3:113:C:C6	2.44	0.51
62:N6:89:LYS:NZ	36:5:375:A:OP2	76.40	0.51
36:5:979:U:H1'	36:5:980:A:N3	2.26	0.51
1:6:1557:U:O2'	1:6:1558:U:H2'	2.10	0.51
14:C2:75:VAL:O	14:C2:79:ALA:N	2.91	0.51
15:C3:26:PHE:HE2	15:C3:66:ILE:HD13	1.75	0.51
15:C3:55:ARG:O	29:D7:47:PHE:HB2	2.34	0.51
17:C5:98:ASN:ND2	17:C5:121:ILE:O	2.37	0.51
1:2:1498:G:H5''	21:C9:72:GLY:HA3	1.92	0.51
2:S0:52:LYS:HG2	23:D1:82:VAL:HG22	4.23	0.51
7:S5:161:ASP:O	30:D8:44:VAL:HA	2.11	0.51
40:L3:275:ARG:NH1	36:5:3045:G:O3'	235.37	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:L5:235:SER:O	42:L5:239:ILE:HG13	2.11	0.51
44:L7:141:TYR:HA	44:L7:144:ILE:HD12	1.92	0.51
47:M0:191:LYS:HE3	47:M0:198:LYS:HD2	1.92	0.51
48:M1:15:GLU:HB3	48:M1:130:VAL:HG13	1.92	0.51
49:M3:174:ARG:CZ	72:O6:9:ILE:HD13	2.41	0.51
36:1:73:C:C2	49:M3:59:ARG:HD3	2.45	0.51
52:M6:31:GLN:HE21	52:M6:32:LYS:N	2.08	0.51
58:N2:41:ILE:HG12	58:N2:79:LEU:HD13	1.91	0.51
63:N7:85:TYR:HE2	63:N7:129:TRP:CE2	3.83	0.51
66:O0:73:GLY:O	66:O0:76:GLU:HG2	2.11	0.51
2:S0:14:ALA:HA	2:S0:17:LEU:HD12	3.02	0.51
5:S3:148:LYS:HB2	35:SM:110:TRP:CZ2	2.45	0.51
6:S4:162:ILE:HG22	6:S4:164:LEU:H	1.75	0.51
6:S4:3:ARG:HG2	1:6:399:A:H4'	319.98	0.51
36:1:160:G:H2'	36:1:161:G:O4'	2.11	0.51
36:1:3165:A:H2'	36:1:3166:C:C6	2.46	0.51
36:1:738:A:H2'	36:1:739:G:C8	2.46	0.51
1:2:1244:A:N3	1:2:1244:A:H3'	2.26	0.51
36:5:1944:U:H2'	36:5:1945:A:H8	1.76	0.51
36:5:2226:U:H2'	36:5:2227:C:C6	2.46	0.51
36:5:2881:C:H2'	36:5:2882:U:H6	1.76	0.51
36:5:3115:C:O2'	36:5:3117:C:N4	2.38	0.51
36:5:437:G:H22	36:5:622:A:H61	1.57	0.51
36:5:763:G:H1	36:5:768:C:H42	1.58	0.51
36:5:93:C:OP2	36:5:2764:C:O2'	2.24	0.51
1:6:1003:A:H1'	1:6:1005:A:N7	2.26	0.51
13:C1:5:LEU:O	13:C1:7:VAL:N	2.37	0.51
19:C7:108:ASP:O	19:C7:112:SER:OG	2.28	0.51
19:C7:57:LEU:O	19:C7:61:ILE:HG13	2.10	0.51
20:C8:54:LEU:H	20:C8:54:LEU:HD22	1.76	0.51
25:D3:6:PRO:HG3	25:D3:14:LYS:HG2	2.11	0.51
31:D9:5:ASN:HB3	31:D9:7:TRP:CZ2	3.59	0.51
39:L2:174:ARG:HA	79:Q3:69:TYR:HE2	2.38	0.51
41:L4:338:LYS:HD2	41:L4:338:LYS:N	2.25	0.51
43:L6:166:LYS:N	43:L6:169:ASP:OD2	2.39	0.51
43:L6:176:PHE:HA	50:M4:114:ASP:HB2	2.20	0.51
48:M1:23:VAL:O	48:M1:25:GLU:N	2.36	0.51
51:M5:109:ARG:O	51:M5:111:ALA:N	2.44	0.51
56:N0:83:SER:OG	56:N0:86:GLY:O	2.46	0.51
56:N0:23:LYS:HD3	57:N1:146:ASN:OD1	7.47	0.51
61:N5:53:HIS:ND1	61:N5:54:TYR:O	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
61:N5:57:LEU:HD11	61:N5:89:LYS:HB2	3.13	0.51
66:O0:13:LYS:HZ1	66:O0:103:THR:HG21	5.62	0.51
68:O2:99:ASN:N	68:O2:99:ASN:OD1	2.52	0.51
69:O3:16:TYR:HB3	69:O3:24:ASN:O	2.62	0.51
73:O7:71:SER:OG	73:O7:72:ARG:N	2.43	0.51
39:L2:171:GLY:O	79:Q3:68:ALA:HB2	2.52	0.51
2:S0:109:ASN:HD21	2:S0:111:ILE:HG22	1.75	0.51
5:S3:61:GLU:O	5:S3:64:ARG:HB2	2.46	0.51
6:S4:240:LYS:HE2	6:S4:240:LYS:H	1.74	0.51
7:S5:99:MET:HG3	7:S5:180:ARG:NH2	3.10	0.51
11:S9:114:TYR:HD2	11:S9:115:LYS:HD2	3.06	0.51
36:1:1014:U:H2'	36:1:1015:U:H5''	1.93	0.51
36:1:1277:C:O2'	36:1:1278:A:H8	1.94	0.51
36:1:2378:C:H2'	36:1:2379:U:H6	1.76	0.51
36:1:789:A:H2'	36:1:790:U:C6	2.46	0.51
1:2:1201:G:H22	1:2:1600:A:H5''	1.76	0.51
1:2:220:A:H5''	1:2:832:U:H1'	1.93	0.51
1:2:702:G:O6	1:2:736:C:N4	2.32	0.51
1:2:783:G:HO2'	1:2:784:C:H6	1.58	0.51
68:O2:105:ARG:NH2	36:5:1412:G:OP1	147.74	0.51
36:5:174:C:H2'	36:5:175:C:H6	1.75	0.51
36:5:438:A:H2'	36:5:494:G:N2	2.26	0.51
1:6:37:U:O2'	1:6:770:A:N1	2.32	0.51
12:C0:24:LYS:HG2	12:C0:29:GLN:HE22	5.35	0.51
12:C0:35:ILE:HG22	12:C0:37:THR:HG23	1.93	0.51
13:C1:22:ASN:ND2	13:C1:24:LYS:HB2	2.25	0.51
20:C8:76:PRO:HB2	20:C8:86:LEU:HD21	1.92	0.51
21:C9:9:VAL:O	21:C9:11:ALA:N	2.43	0.51
39:L2:10:LYS:HA	39:L2:16:PHE:CD2	2.46	0.51
41:L4:359:LEU:HD13	56:N0:8:GLN:NE2	2.26	0.51
48:M1:108:GLU:HB3	48:M1:110:ILE:HG12	1.91	0.51
52:M6:108:ILE:HG12	52:M6:160:ARG:HE	6.02	0.51
78:Q2:6:LYS:HG3	78:Q2:93:LEU:O	2.11	0.51
2:S0:163:ASN:HD21	2:S0:165:ARG:HG2	1.76	0.51
11:S9:106:GLU:O	11:S9:111:THR:OG1	3.64	0.51
34:SR:106:HIS:CD2	34:SR:110:VAL:HG22	2.48	0.51
36:1:2709:C:H2'	36:1:2710:C:H6	1.76	0.51
1:2:791:A:H2'	1:2:792:U:H6	1.76	0.51
36:5:1190:A:H4'	36:5:1191:U:OP1	2.10	0.51
36:5:1692:U:C4	36:5:1693:C:N4	2.79	0.51
36:5:252:U:H4'	36:5:253:A:C5'	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:722:G:C6	36:5:749:C:N3	2.78	0.51
1:6:1142:A:H2'	1:6:1143:A:O4'	2.10	0.51
1:6:751:G:H2'	1:6:752:A:H8	1.76	0.51
16:C4:81:VAL:HG13	16:C4:115:ILE:HG23	4.45	0.51
17:C5:44:ARG:HG3	17:C5:84:ILE:HD11	1.92	0.51
18:C6:7:VAL:HG12	18:C6:8:GLN:H	4.18	0.51
26:D4:60:PHE:O	1:6:523:G:H5'	413.95	0.51
39:L2:95:SER:OG	39:L2:100:ASN:ND2	2.43	0.51
36:1:2242:A:H5'	39:L2:243:THR:O	2.11	0.51
39:L2:47:GLN:HG2	39:L2:60:LYS:HB2	4.55	0.51
40:L3:320:ASP:N	40:L3:320:ASP:OD2	2.44	0.51
41:L4:257:LYS:HA	41:L4:260:GLN:HG2	1.92	0.51
41:L4:80:GLY:HA2	41:L4:85:SER:CB	2.41	0.51
42:L5:64:ILE:HG22	42:L5:75:LEU:HB3	1.92	0.51
43:L6:40:LEU:HD13	43:L6:84:VAL:HG21	1.93	0.51
36:1:770:G:OP2	49:M3:171:ARG:NH2	2.44	0.51
53:M7:69:ARG:HG2	53:M7:79:THR:HB	2.41	0.51
56:N0:12:ARG:HH12	56:N0:15:PRO:HG3	1.76	0.51
56:N0:1:MET:HB2	56:N0:118:PHE:CD1	2.46	0.51
63:N7:20:GLY:HA3	63:N7:136:PHE:HE1	3.35	0.51
63:N7:81:LEU:HD11	70:O4:90:ILE:HG23	1.93	0.51
64:N8:6:THR:HG23	64:N8:8:THR:H	1.76	0.51
68:O2:124:GLY:O	68:O2:126:LEU:N	2.97	0.51
73:O7:17:THR:HG22	73:O7:18:LEU:H	1.75	0.51
2:S0:148:ASP:OD2	2:S0:165:ARG:NH2	2.44	0.51
5:S3:132:LYS:HE3	5:S3:192:PRO:HD2	1.93	0.51
36:1:158:G:H2'	36:1:159:A:C8	2.46	0.51
36:1:289:A:H5''	51:M5:97:SER:HB3	1.93	0.51
36:1:540:U:H2'	36:1:541:U:O4'	2.11	0.51
36:1:616:G:H2'	36:1:617:G:H8	1.75	0.51
36:1:956:U:H2'	36:1:957:C:C6	2.46	0.51
1:2:1563:C:H2'	1:2:1564:U:C6	2.46	0.51
1:2:252:U:H2'	1:2:253:A:H8	1.76	0.51
1:2:681:U:H2'	1:2:682:C:C6	2.46	0.51
1:2:688:G:H2'	1:2:689:G:H8	1.75	0.51
1:2:87:C:H1'	1:2:168:A:N1	2.26	0.51
36:5:2116:G:OP1	36:5:2118:C:N4	2.43	0.51
36:5:2651:G:H4'	36:5:2652:U:OP2	2.11	0.51
36:5:495:G:H2'	36:5:496:C:O4'	2.10	0.51
36:5:528:U:H2'	36:5:529:A:H8	1.76	0.51
36:5:601:U:O4	36:5:602:A:N6	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D6:76:SER:HG	1:6:1793:G:H22	320.80	0.51
1:6:493:U:H2'	1:6:494:U:H5''	1.92	0.51
11:S9:143:ILE:HG21	1:6:768:C:H1'	419.54	0.51
71:O5:7:TYR:HE2	38:8:86:U:H2'	19.88	0.51
17:C5:16:SER:HA	17:C5:21:ASP:HA	1.92	0.51
2:S0:55:GLU:HB3	23:D1:79:LEU:HD11	1.92	0.51
25:D3:52:ILE:HA	25:D3:101:GLU:HA	2.82	0.51
39:L2:70:ARG:HH21	39:L2:72:ARG:NH2	6.82	0.51
42:L5:269:SER:HB2	37:7:1:G:H21	316.67	0.51
44:L7:143:THR:HG22	44:L7:241:LYS:HE3	1.92	0.51
45:L8:133:LYS:HD2	45:L8:138:HIS:HE1	1.76	0.51
52:M6:85:ARG:O	52:M6:87:MET:N	2.44	0.51
54:M8:67:ILE:HG23	54:M8:81:VAL:HG11	3.17	0.51
60:N4:46:PRO:O	60:N4:52:THR:OG1	2.64	0.51
61:N5:58:ASP:OD1	71:O5:25:LYS:NZ	2.37	0.51
62:N6:60:ARG:HB2	62:N6:103:LYS:HB3	1.92	0.51
66:O0:55:GLU:HB2	70:O4:94:LEU:HD11	2.78	0.51
36:1:1367:G:OP1	68:O2:45:ARG:NH2	2.44	0.51
68:O2:75:LEU:HD23	68:O2:95:GLU:HB3	1.92	0.51
70:O4:58:ARG:HG3	70:O4:59:PRO:HD2	2.60	0.51
70:O4:89:ILE:HG22	70:O4:90:ILE:HD13	1.93	0.51
72:O6:66:GLU:OE1	72:O6:70:ARG:NH2	2.44	0.51
2:S0:77:SER:OG	2:S0:78:SER:N	2.43	0.51
2:S0:7:PHE:HZ	23:D1:43:GLY:HA2	2.70	0.51
3:S1:143:THR:O	3:S1:208:GLN:NE2	2.44	0.51
4:S2:69:ILE:HD13	4:S2:136:VAL:HG21	2.30	0.51
10:S8:38:ILE:HG12	10:S8:96:LEU:HD11	1.92	0.51
11:S9:64:GLU:HG2	11:S9:69:ARG:NH1	2.26	0.51
34:SR:184:ASN:N	34:SR:184:ASN:OD1	2.43	0.51
34:SR:267:PRO:HD2	34:SR:269:TYR:CE1	3.74	0.51
36:1:1954:G:H5'	36:1:1955:U:OP2	2.10	0.50
36:1:71:A:C2	36:1:2778:G:H1'	2.46	0.50
36:1:3283:U:H2'	36:1:3284:G:H8	1.77	0.50
36:1:3325:G:H5'	67:O1:104:LEU:O	2.12	0.50
36:1:789:A:H2'	36:1:790:U:H6	1.76	0.50
1:2:407:A:O2'	1:2:1671:A:N3	2.40	0.50
1:2:1789:G:OP2	16:C4:132:ARG:NH2	2.42	0.50
1:2:324:U:OP1	13:C1:133:LYS:NZ	2.27	0.50
1:2:417:A:H4'	1:2:418:G:O5'	2.11	0.50
36:5:2389:C:H42	36:5:2990:G:H1	1.59	0.50
1:6:93:A:C6	1:6:398:G:C6	3.00	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:628:G:N1	1:6:970:A:OP2	2.44	0.50
1:2:1566:U:H4'	20:C8:37:GLY:O	2.11	0.50
22:D0:24:ILE:HG12	22:D0:116:VAL:HG13	1.91	0.50
24:D2:122:SER:OG	24:D2:123:GLY:N	2.44	0.50
26:D4:12:VAL:HG22	26:D4:23:PHE:HB3	3.23	0.50
39:L2:107:VAL:HB	39:L2:111:THR:OG1	2.11	0.50
40:L3:56:ILE:HD11	40:L3:359:ILE:HG12	2.01	0.50
42:L5:155:THR:HA	42:L5:179:ARG:HA	2.08	0.50
45:L8:130:TYR:CE1	45:L8:202:GLU:HB3	2.46	0.50
49:M3:57:VAL:HG22	49:M3:147:ILE:HD13	3.76	0.50
53:M7:29:THR:O	53:M7:32:THR:HG23	5.56	0.50
59:N3:120:LYS:HB2	59:N3:137:VAL:CG2	2.41	0.50
61:N5:58:ASP:O	61:N5:62:VAL:HG23	2.39	0.50
67:O1:72:ARG:HD3	67:O1:104:LEU:HD12	2.89	0.50
39:L2:57:PRO:HD3	79:Q3:53:GLY:HA3	3.04	0.50
2:S0:36:TYR:OH	23:D1:70:ASN:ND2	2.41	0.50
4:S2:41:LEU:HD11	4:S2:56:ILE:HD13	2.86	0.50
6:S4:26:CYS:SG	1:6:461:G:H5''	363.73	0.50
6:S4:65:LEU:HD13	6:S4:80:THR:HA	3.58	0.50
36:1:1355:A:H5'	36:1:1357:G:H1'	1.92	0.50
36:1:1582:C:HO2'	36:1:1583:A:P	2.34	0.50
36:1:3228:C:H4'	36:1:3229:G:O5'	2.11	0.50
1:2:1433:G:C4	31:D9:41:GLN:HB3	2.45	0.50
1:2:1660:A:H2'	1:2:1661:U:C6	2.47	0.50
1:2:415:C:O2'	1:2:418:G:O6	2.22	0.50
1:2:685:A:O2'	1:2:686:C:OP1	2.27	0.50
1:2:891:A:H2'	1:2:892:A:C8	2.46	0.50
1:2:904:G:H2'	1:2:905:A:C8	2.46	0.50
36:5:3153:U:H4'	36:5:3154:C:H5'	1.92	0.50
36:5:748:U:H2'	36:5:749:C:C6	2.46	0.50
36:5:899:U:H2'	36:5:900:G:H8	1.76	0.50
1:6:188:A:H2'	1:6:189:C:O4'	2.10	0.50
27:D5:71:ILE:HG21	27:D5:76:ALA:HB2	4.74	0.50
28:D6:84:VAL:HG13	28:D6:85:ARG:H	1.77	0.50
40:L3:186:GLY:O	40:L3:190:GLU:HB2	4.01	0.50
40:L3:346:THR:HG22	40:L3:351:LEU:HD11	3.68	0.50
41:L4:138:ARG:HH21	41:L4:240:PRO:HB2	1.77	0.50
47:M0:54:SER:HB3	47:M0:135:ILE:HD11	1.92	0.50
49:M3:94:GLY:HA3	71:O5:116:TYR:OH	2.12	0.50
52:M6:158:ALA:O	52:M6:162:VAL:HG23	2.11	0.50
56:N0:12:ARG:HB3	56:N0:24:LEU:HD23	2.05	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:N1:158:THR:OG1	57:N1:159:PHE:N	2.43	0.50
60:N4:25:ASP:O	60:N4:26:SER:HB2	2.30	0.50
64:N8:118:ILE:HD13	64:N8:118:ILE:H	1.75	0.50
54:M8:175:ALA:HB2	64:N8:56:VAL:HG22	1.92	0.50
66:O0:24:THR:HG23	66:O0:30:THR:HG22	2.50	0.50
67:O1:29:ALA:HB3	67:O1:30:PRO:HD3	1.93	0.50
69:O3:26:ASN:OD1	69:O3:88:ASN:ND2	2.42	0.50
72:O6:55:ARG:O	72:O6:58:ILE:HG12	2.11	0.50
6:S4:28:ALA:O	1:6:448:C:H4'	365.12	0.50
7:S5:162:VAL:HG21	7:S5:166:ARG:HH11	1.76	0.50
9:S7:124:LYS:HA	9:S7:127:GLU:HB2	2.67	0.50
36:1:1103:A:N6	36:1:1363:A:H1'	2.26	0.50
36:1:158:G:H2'	36:1:159:A:H8	1.76	0.50
36:1:2897:A:H2'	36:1:2899:C:H5''	1.93	0.50
36:1:2904:U:H2'	36:1:2905:U:C6	2.46	0.50
36:1:3283:U:H2'	36:1:3284:G:C8	2.46	0.50
1:2:924:A:H2'	1:2:925:G:C8	2.46	0.50
36:5:1784:G:H2'	36:5:1785:U:O4'	2.11	0.50
36:5:2572:C:H2'	36:5:2572:C:OP2	2.12	0.50
36:5:916:G:H5'	36:5:917:A:OP1	2.10	0.50
1:6:1372:U:H2'	1:6:1373:C:C6	2.46	0.50
7:S5:25:LEU:HD12	18:C6:61:SER:HA	3.57	0.50
20:C8:145:ARG:HG3	35:SM:68:ARG:NH1	6.36	0.50
28:D6:35:ALA:HB3	28:D6:37:LYS:HE3	1.92	0.50
33:E1:139:LEU:HB2	33:E1:150:VAL:O	2.11	0.50
42:L5:36:LEU:HD23	36:5:2748:A:N3	254.85	0.50
44:L7:173:LEU:HD21	44:L7:198:ALA:HA	2.53	0.50
44:L7:208:SER:OG	44:L7:209:ASN:N	2.84	0.50
44:L7:228:SER:HA	44:L7:232:ARG:HH21	2.62	0.50
44:L7:52:GLN:O	44:L7:56:GLU:HG2	2.10	0.50
46:L9:47:LYS:HB2	50:M4:7:VAL:HB	1.94	0.50
46:L9:70:THR:O	46:L9:74:LEU:HG	2.46	0.50
47:M0:48:LEU:O	47:M0:139:ARG:HA	2.35	0.50
47:M0:49:CYS:HB3	47:M0:168:SER:HB3	1.94	0.50
52:M6:12:LYS:HG3	52:M6:40:GLU:HB3	2.52	0.50
68:O2:24:ARG:HG2	68:O2:25:TYR:CZ	3.09	0.50
71:O5:31:LEU:HD13	71:O5:44:ILE:HA	1.92	0.50
71:O5:49:LYS:HE2	38:8:63:G:O3'	49.61	0.50
5:S3:125:TYR:HE1	35:SM:134:ASP:OD2	1.94	0.50
5:S3:170:THR:HG22	5:S3:187:LYS:HG3	5.52	0.50
6:S4:64:ILE:HD11	26:D4:18:LEU:HG	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:S8:184:LEU:HD23	10:S8:188:GLU:HG2	4.50	0.50
1:2:257:A:O2'	10:S8:64:ASN:OD1	2.26	0.50
20:C8:125:ILE:HD11	35:SM:57:ASN:HB3	1.93	0.50
36:1:1308:A:N1	36:1:2381:G:O2'	2.31	0.50
36:1:1638:A:N1	36:1:1736:G:O2'	2.36	0.50
36:1:2445:A:N6	36:1:2502:A:N1	2.60	0.50
36:1:3096:C:H2'	36:1:3097:C:C6	2.46	0.50
36:1:760:G:H1'	36:1:770:G:N2	2.26	0.50
1:2:1756[A]:A:H2'	1:2:1757:G:H8	1.77	0.50
1:2:322:G:O2'	10:S8:10:LYS:NZ	2.42	0.50
36:5:2254:U:H2'	36:5:2261:G:H22	1.76	0.50
52:M6:68:ARG:NH1	36:5:2988:C:OP1	217.86	0.50
36:5:789:A:H2'	36:5:790:U:C6	2.46	0.50
1:6:1159:C:H5''	1:6:1160:A:H5'	1.93	0.50
1:6:94:U:H2'	1:6:95:G:O4'	2.12	0.50
13:C1:75:VAL:HG12	13:C1:119:VAL:HA	1.91	0.50
15:C3:26:PHE:CE2	15:C3:66:ILE:HD13	2.46	0.50
21:C9:27:LYS:HB3	21:C9:111:ILE:HD11	1.93	0.50
41:L4:358:THR:HG21	57:N1:148:PRO:HD2	1.94	0.50
45:L8:68:ARG:HA	45:L8:236:GLY:O	4.18	0.50
46:L9:44:THR:HG22	36:5:3186:A:H1'	325.47	0.50
47:M0:89:VAL:HG13	47:M0:136:PHE:HE1	1.75	0.50
49:M3:161:ASP:OD1	49:M3:162:ASN:N	3.67	0.50
49:M3:73:ARG:HD2	36:5:76:G:H3'	82.92	0.50
51:M5:113:LEU:HB2	51:M5:134:LEU:HD23	4.20	0.50
53:M7:29:THR:HG22	53:M7:87:SER:OG	2.10	0.50
36:1:670:C:OP1	54:M8:147:ARG:NH2	2.44	0.50
46:L9:4:ILE:HG22	56:N0:142:GLN:OE1	2.76	0.50
62:N6:111:LEU:HD23	62:N6:116:LYS:HG2	1.93	0.50
63:N7:83:THR:HG23	63:N7:85:TYR:N	2.23	0.50
72:O6:36:ARG:HE	72:O6:40:VAL:HG21	3.82	0.50
73:O7:64:MET:O	73:O7:68:LYS:HB2	2.12	0.50
6:S4:122:LYS:HE3	6:S4:164:LEU:HD21	1.92	0.50
9:S7:110:GLN:HE21	9:S7:110:GLN:HA	1.77	0.50
36:1:1658:G:H2'	36:1:1659:U:C6	2.47	0.50
36:1:177:U:H2'	36:1:178:U:O4'	2.12	0.50
36:1:2233:A:H2'	36:1:2234:G:O4'	2.12	0.50
36:1:679:U:H1'	36:1:788:C:H1'	1.92	0.50
36:5:3349:C:H42	36:5:3356:G:H1	1.59	0.50
36:5:736:A:C5	36:5:737:G:H1'	2.46	0.50
42:L5:260:PHE:HE2	37:7:121:U:H5'	320.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
73:O7:63:ARG:NH2	38:8:58:G:O6	79.08	0.50
16:C4:50:ALA:HB3	16:C4:53:ASP:HB2	1.92	0.50
17:C5:121:ILE:HG23	17:C5:123:TYR:H	1.77	0.50
17:C5:21:ASP:HB3	17:C5:24:LYS:NZ	2.27	0.50
18:C6:131:GLY:HA3	18:C6:136:SER:O	2.60	0.50
24:D2:15:ASN:HD21	24:D2:71:LYS:HG3	1.77	0.50
39:L2:204:MET:HE2	39:L2:209:HIS:HB2	1.92	0.50
41:L4:327:LEU:HA	44:L7:166:ASN:HD21	1.75	0.50
45:L8:89:GLU:HA	45:L8:92:LYS:HB2	1.94	0.50
45:L8:97:TYR:O	45:L8:132:VAL:HG12	2.12	0.50
46:L9:23:ARG:NH2	46:L9:42:ASP:OD2	2.45	0.50
48:M1:23:VAL:HG13	48:M1:29:ARG:HH11	1.76	0.50
53:M7:30:ARG:HA	53:M7:119:VAL:HG11	2.19	0.50
36:1:943:U:H3'	64:N8:13:GLY:HA2	1.94	0.50
49:M3:64:LYS:HG3	64:N8:69:TRP:CG	2.46	0.50
71:O5:111:PHE:HZ	36:5:256:G:H4'	53.07	0.50
1:2:1514:U:O2'	5:S3:5:ILE:O	2.30	0.50
6:S4:101:LEU:HB3	6:S4:109:PHE:HE1	1.76	0.50
6:S4:18:TRP:HH2	6:S4:31:PRO:HD3	2.19	0.50
9:S7:155:ASP:OD2	9:S7:156:SER:N	2.35	0.50
36:1:1393:A:N3	36:1:1419:A:O2'	2.42	0.50
36:1:201:A:H2'	36:1:202:G:C8	2.46	0.50
36:1:3150:A:H2'	36:1:3151:U:O4'	2.12	0.50
36:1:379:C:H2'	36:1:380:U:H6	1.77	0.50
36:1:63:A:N3	36:1:78:U:O2'	2.39	0.50
36:5:1152:G:H22	36:5:1200:A:H61	1.57	0.50
36:5:1899:G:H22	36:5:2335:G:H5'	1.76	0.50
40:L3:242:THR:HG22	36:5:2948:C:O2'	214.97	0.50
41:L4:93:MET:HB2	36:5:658:G:N2	145.81	0.50
1:6:1304:G:H5'	1:6:1322:A:OP2	2.12	0.50
1:6:1578:U:H2'	1:6:1579:U:H6	1.76	0.50
1:6:1698:G:N2	1:6:1699:G:N3	2.59	0.50
1:6:968:U:OP1	1:6:1033:C:O2'	2.28	0.50
14:C2:108:UNK:O	14:C2:110:UNK:N	2.71	0.50
25:D3:69:ARG:HD3	25:D3:117:ILE:HG23	1.92	0.50
26:D4:112:LYS:HE3	26:D4:116:LYS:HD2	1.94	0.50
27:D5:85:LYS:HG3	27:D5:86:GLU:N	2.27	0.50
29:D7:14:SER:HB2	29:D7:17:ARG:HE	1.77	0.50
39:L2:189:TYR:HA	39:L2:192:LYS:HB2	1.92	0.50
41:L4:138:ARG:HD3	41:L4:140:HIS:CD2	2.87	0.50
43:L6:28:GLN:OE1	43:L6:61:ASN:ND2	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:L6:40:LEU:HB3	43:L6:84:VAL:CG1	4.18	0.50
46:L9:12:VAL:HG13	46:L9:16:VAL:HG23	1.94	0.50
36:1:3115:C:OP1	46:L9:62:ARG:NH2	2.44	0.50
49:M3:64:LYS:HE3	64:N8:69:TRP:CD1	2.46	0.50
50:M4:116:GLU:O	50:M4:120:VAL:HG23	2.11	0.50
52:M6:55:HIS:HA	52:M6:58:LEU:HB2	1.93	0.50
36:1:784:A:C8	54:M8:69:ARG:HG3	2.46	0.50
68:O2:119:VAL:HG12	68:O2:122:PRO:HD3	1.93	0.50
69:O3:49:ILE:HA	69:O3:99:ARG:O	2.39	0.50
74:O8:58:ASP:HB3	74:O8:61:LYS:HB2	1.93	0.50
6:S4:51:ARG:O	6:S4:52:LEU:HD23	3.58	0.50
36:1:1622:U:H2'	36:1:1623:G:H8	1.77	0.50
36:1:1805:C:H2'	36:1:1806:A:C8	2.46	0.50
36:1:3182:G:H2'	36:1:3183:A:O4'	2.11	0.50
36:1:739:G:H2'	36:1:740:G:H8	1.75	0.50
1:2:1091:A:C8	1:2:1092:A:C6	3.00	0.50
1:2:1300:A:OP1	4:S2:99:LYS:NZ	2.36	0.50
1:2:1583:A:N6	1:2:1612:U:OP2	2.39	0.50
1:2:542:A:H8	1:2:543:C:H5'	1.76	0.50
1:2:926:A:H1'	1:2:988:A:C2	2.46	0.50
36:5:1039:U:H2'	36:5:1040:A:C8	2.47	0.50
36:5:209:A:H1'	36:5:212:G:N2	2.27	0.50
1:6:1524:A:H2'	1:6:1525:A:C8	2.47	0.50
77:Q1:21:ARG:HH11	1:6:1654:G:P	281.21	0.50
1:6:570:A:H8	1:6:570:A:OP2	1.95	0.50
15:C3:136:PRO:HD2	15:C3:139:TRP:HD1	3.51	0.50
22:D0:80:GLU:OE1	31:D9:44:ARG:NH1	2.51	0.50
15:C3:18:TYR:HA	24:D2:56:HIS:CD2	4.95	0.50
25:D3:43:PHE:CD1	25:D3:49:ALA:HB3	2.66	0.50
41:L4:293:SER:O	41:L4:297:SER:OG	2.16	0.50
49:M3:192:GLU:O	49:M3:194:GLU:N	3.06	0.50
36:1:304:G:C6	51:M5:178:HIS:HD2	2.29	0.50
53:M7:22:LEU:HD12	53:M7:146:ILE:HG13	2.36	0.50
53:M7:26:PHE:HZ	36:5:412:G:H5'	150.84	0.50
53:M7:24:VAL:HB	53:M7:29:THR:HG21	1.91	0.50
54:M8:37:ALA:O	54:M8:46:LYS:HE3	2.83	0.50
57:N1:105:PHE:O	57:N1:109:VAL:HG23	2.36	0.50
58:N2:14:THR:HG23	58:N2:66:VAL:HG13	1.94	0.50
36:1:2916:U:C1'	59:N3:44:SER:HB3	2.41	0.50
59:N3:89:ASP:OD1	59:N3:91:VAL:HG12	5.34	0.50
68:O2:11:LYS:O	68:O2:13:HIS:N	2.34	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
70:O4:41:ARG:HD2	70:O4:56:THR:HG21	3.92	0.50
73:O7:27:PHE:HA	73:O7:34:CYS:HA	1.92	0.50
3:S1:142:PHE:O	3:S1:207:LEU:HA	2.64	0.50
5:S3:22:ASN:O	5:S3:26:THR:OG1	2.21	0.50
8:S6:30:LYS:HE3	8:S6:34:GLN:HG2	3.53	0.50
11:S9:74:ASN:O	11:S9:78:ARG:HB3	3.03	0.50
35:SM:70:ASN:OD1	35:SM:70:ASN:N	2.43	0.50
34:SR:16:HIS:CD2	34:SR:37:SER:HB2	3.30	0.50
36:1:1814:A:H1'	36:1:1815:U:C2	2.47	0.50
36:1:2152:A:H2'	36:1:2153:U:H6	1.76	0.50
36:1:304:G:C5	51:M5:178:HIS:HD2	2.30	0.50
36:1:738:A:H2'	36:1:739:G:H8	1.76	0.50
1:2:1251:U:H5'	33:E1:135:HIS:CD2	2.43	0.50
1:2:1316:G:P	19:C7:7:LYS:HZ3	2.35	0.50
1:2:1489:U:OP2	5:S3:9:ARG:NH2	2.42	0.50
1:2:393:C:H2'	1:2:394:C:C6	2.46	0.50
62:N6:19:TYR:CZ	36:5:216:G:H4'	72.44	0.50
36:5:235:A:H2'	36:5:236:G:C8	2.47	0.50
51:M5:44:ARG:HH22	36:5:269:G:P	125.72	0.50
36:5:874:U:H5''	36:5:2950:G:OP1	2.11	0.50
69:O3:56:SER:OG	36:5:3170:A:OP2	203.92	0.50
41:L4:271:LYS:NZ	36:5:695:C:OP1	103.60	0.50
1:6:1237:G:H2'	1:6:1238:A:C8	2.47	0.50
1:6:1357:A:H2'	1:6:1358:G:C8	2.47	0.50
1:6:766:U:H5'	1:6:767:U:H5''	1.93	0.50
42:L5:16:PHE:O	37:7:11:A:N6	294.77	0.50
36:1:2341:A:P	40:L3:247:ARG:HH22	2.35	0.50
45:L8:65:LEU:HD12	51:M5:25:VAL:HG13	1.94	0.50
52:M6:67:THR:O	52:M6:71:PHE:HE1	2.61	0.50
59:N3:74:MET:HG3	59:N3:102:ILE:HG23	5.83	0.50
65:N9:38:LYS:HB3	65:N9:41:ARG:NH1	4.98	0.50
2:S0:179:ARG:HD3	2:S0:183:ARG:HH11	1.77	0.50
2:S0:30:GLN:OE1	2:S0:31:VAL:N	5.96	0.50
4:S2:101:VAL:HG13	4:S2:115:ILE:HG12	1.94	0.50
5:S3:107:PHE:O	5:S3:111:ASN:N	3.99	0.50
10:S8:3:ILE:O	10:S8:30:GLY:N	2.44	0.50
10:S8:6:ASP:HB3	10:S8:28:GLU:OE2	2.34	0.50
36:1:2424:A:C2	39:L2:230:VAL:HG21	2.47	0.50
36:1:627:U:H2'	36:1:628:A:C8	2.47	0.50
1:2:1049:U:H2'	1:2:1050:G:C8	2.46	0.50
1:2:894:U:H2'	1:2:895:G:C8	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:3:115:G:H2'	37:3:116:C:C6	2.47	0.50
58:N2:42:LYS:NZ	36:5:1686:U:OP1	177.34	0.50
39:L2:69:TYR:OH	36:5:2557:A:OP1	191.42	0.50
39:L2:2:GLY:N	36:5:2608:G:OP1	183.37	0.50
36:5:3288:G:C4	36:5:3289:G:C8	3.00	0.50
36:5:549:U:H2'	36:5:550:A:C8	2.47	0.50
69:O3:21:ARG:O	36:5:634:C:H5'	223.13	0.50
1:6:296:U:H2'	1:6:297:U:O4'	2.11	0.50
61:N5:49:LYS:NZ	38:8:135:G:OP1	78.38	0.50
1:2:866:G:H5''	15:C3:3:ARG:H	1.75	0.50
15:C3:56:ASP:OD1	29:D7:52:THR:OG1	2.23	0.50
16:C4:31:THR:OG1	16:C4:32:ASP:N	2.45	0.50
18:C6:114:ARG:O	18:C6:115:THR:HB	3.88	0.50
25:D3:23:ARG:HD2	25:D3:26:GLU:OE1	2.87	0.50
32:E0:50:VAL:O	32:E0:52:GLY:N	2.45	0.50
33:E1:86:THR:O	33:E1:87:THR:OG1	2.67	0.50
39:L2:179:LEU:O	39:L2:181:LYS:N	2.44	0.50
40:L3:126:LYS:HD2	36:5:3294:A:H5''	191.31	0.50
47:M0:56:GLU:HB2	47:M0:58:GLU:OE1	3.37	0.50
52:M6:8:VAL:HG13	52:M6:34:VAL:HG23	1.93	0.50
53:M7:19:GLY:HA3	53:M7:22:LEU:HD11	2.32	0.50
57:N1:79:MET:HB2	57:N1:84:TYR:CE2	3.50	0.50
58:N2:35:LYS:NZ	58:N2:39:ASP:OD1	2.25	0.50
62:N6:45:ILE:HD13	62:N6:122:LYS:HE3	4.77	0.50
71:O5:83:LYS:O	71:O5:89:ARG:NE	2.94	0.50
49:M3:180:ARG:HD3	72:O6:11:LEU:HD11	2.02	0.50
3:S1:109:LYS:HG3	3:S1:113:MET:HE3	1.94	0.50
3:S1:21:VAL:HG23	3:S1:22:ASP:H	1.90	0.50
6:S4:103:TYR:HE1	6:S4:109:PHE:CE2	2.29	0.50
34:SR:84:SER:OG	34:SR:85:TRP:N	2.48	0.50
36:1:1313:G:O2'	36:1:1318:A:N1	2.42	0.49
36:1:2100:A:N7	36:1:2101:C:N4	2.60	0.49
36:1:2882:U:H2'	36:1:2883:U:C6	2.47	0.49
36:1:821:U:H2'	36:1:822:G:H8	1.77	0.49
36:1:873:C:H2'	36:1:875:G:O4'	2.12	0.49
1:2:1483:A:H2'	1:2:1484:G:C8	2.47	0.49
1:2:1552:U:H2'	1:2:1553:G:O4'	2.13	0.49
1:2:552:G:C6	1:2:553:G:C6	3.00	0.49
1:2:704:C:N4	1:2:734:A:O2'	2.30	0.49
1:2:97:C:H2'	1:2:98:U:C6	2.47	0.49
1:2:980:G:OP2	1:2:1014:G:O2'	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:1840:U:H4'	36:5:1841:A:H5'	1.93	0.49
36:5:496:C:N3	36:5:617:G:N1	2.59	0.49
36:5:602:A:H2'	36:5:603:A:C8	2.47	0.49
68:O2:55:ILE:HB	36:5:947:G:H5'	187.29	0.49
33:E1:83:LYS:NZ	1:6:1211:A:OP1	367.91	0.49
1:6:1382:A:O2'	1:6:1383:G:H5''	2.11	0.49
1:6:1511:U:H2'	1:6:1512:G:C8	2.47	0.49
1:6:1466:G:O2'	1:6:1602:C:OP1	2.28	0.49
1:6:1672:G:H2'	1:6:1673:G:C8	2.47	0.49
1:6:675:U:H2'	1:6:676:G:C8	2.46	0.49
12:C0:81:ASN:ND2	12:C0:81:ASN:O	2.45	0.49
13:C1:108:PRO:HG3	13:C1:134:THR:HB	3.43	0.49
13:C1:135:VAL:O	13:C1:136:ARG:NH1	2.42	0.49
7:S5:112:ARG:NH1	18:C6:43:ILE:HD11	2.27	0.49
28:D6:53:LEU:HD11	28:D6:62:TYR:CD2	2.47	0.49
39:L2:5:ILE:HG22	39:L2:208:ASP:O	2.12	0.49
40:L3:370:PHE:HD1	40:L3:375:GLU:HG2	1.76	0.49
42:L5:160:PHE:CE2	42:L5:179:ARG:HB3	2.76	0.49
43:L6:43:LEU:HD11	43:L6:85:ILE:HG13	2.12	0.49
46:L9:166:ARG:HD2	46:L9:168:ARG:HH11	14.57	0.49
51:M5:31:ARG:HG3	51:M5:129:TYR:OH	3.72	0.49
57:N1:38:ASP:O	57:N1:64:VAL:HG23	2.12	0.49
69:O3:49:ILE:HG23	69:O3:100:ILE:HG13	1.92	0.49
73:O7:58:THR:O	73:O7:61:THR:OG1	3.42	0.49
75:O9:30:ARG:HB2	75:O9:33:ASN:HB2	1.93	0.49
76:Q0:98:LYS:HG3	76:Q0:118:THR:HG21	4.30	0.49
2:S0:71:GLU:O	2:S0:96:THR:HG22	2.14	0.49
4:S2:99:LYS:HG3	4:S2:117:THR:HG22	5.61	0.49
8:S6:67:VAL:HG23	8:S6:68:LEU:O	2.30	0.49
10:S8:105:ASP:O	10:S8:108:PRO:HD2	4.87	0.49
10:S8:57:ALA:HB2	10:S8:177:GLY:HA2	2.28	0.49
36:1:1550:C:H2'	36:1:1551:C:C6	2.47	0.49
36:1:1562:C:O2'	36:1:1563:C:O5'	2.30	0.49
36:1:1764:U:C5	36:1:1765:U:H1'	2.47	0.49
36:1:169:U:H3	36:1:253:A:H61	1.59	0.49
36:1:3016:A:H2'	36:1:3017:A:C8	2.47	0.49
36:1:796:U:H2'	36:1:797:U:C6	2.44	0.49
1:2:1349:G:H2'	1:2:1350:U:C6	2.48	0.49
1:2:1218:G:N2	1:2:1444:A:OP2	2.35	0.49
1:2:1494:C:H2'	1:2:1495:C:H6	1.77	0.49
1:2:1553:G:H22	1:2:1555:A:H3'	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:852:C:OP1	55:M9:172:ARG:HB3	2.11	0.49
37:3:118:A:H2'	37:3:119:U:O4'	2.12	0.49
38:4:37:A:H5''	38:4:39:G:O4'	2.12	0.49
1:6:1229:G:H22	1:6:1255:G:H2'	1.77	0.49
1:6:413:U:H2'	1:6:414:C:C6	2.47	0.49
42:L5:155:THR:HG23	37:7:36:C:H4'	270.08	0.49
18:C6:112:TYR:O	18:C6:114:ARG:HG2	7.02	0.49
1:2:1390:U:C6	19:C7:3:ARG:HG2	2.48	0.49
20:C8:24:GLY:HA2	20:C8:58:ALA:HB3	2.44	0.49
22:D0:106:ILE:HG13	22:D0:107:THR:H	1.77	0.49
2:S0:185:ARG:HG3	23:D1:47:PRO:HD3	1.94	0.49
39:L2:250:GLN:HG2	39:L2:251:LYS:H	4.29	0.49
40:L3:14:LEU:O	40:L3:17:LEU:HD22	2.12	0.49
42:L5:270:LYS:HG3	42:L5:273:ARG:HD2	1.94	0.49
44:L7:54:GLU:OE1	44:L7:186:HIS:NE2	2.73	0.49
47:M0:81:GLY:O	47:M0:84:ALA:N	3.52	0.49
54:M8:151:ARG:HD2	36:5:781:G:OP1	160.85	0.49
44:L7:110:ARG:CZ	54:M8:3:ILE:HD12	4.64	0.49
57:N1:100:LYS:HB3	36:5:990:U:H4'	258.78	0.49
59:N3:85:TRP:CE2	59:N3:93:LEU:HD21	2.47	0.49
78:Q2:10:THR:HA	78:Q2:20:HIS:HD2	2.91	0.49
3:S1:202:LYS:O	3:S1:202:LYS:HD3	5.46	0.49
5:S3:133:GLY:HA2	5:S3:155:GLY:HA3	2.62	0.49
1:2:448:C:OP1	6:S4:29:PRO:HD3	2.12	0.49
9:S7:126:LEU:HD21	9:S7:152:VAL:HG21	5.19	0.49
11:S9:42:ILE:O	11:S9:46:SER:OG	2.35	0.49
36:1:1233:G:H22	36:1:1255:C:N4	2.10	0.49
36:1:1782:U:H2'	36:1:1783:U:O4'	2.11	0.49
36:1:2601:A:H2'	36:1:2602:G:C8	2.47	0.49
36:1:2853:A:H4'	47:M0:63:GLU:O	2.11	0.49
36:1:3251:U:H2'	36:1:3252:G:O4'	2.12	0.49
36:1:3343:G:N2	36:1:3361:G:H2'	2.28	0.49
36:1:38:U:H2'	36:1:39:A:O4'	2.12	0.49
1:2:870:C:H2'	1:2:871:G:H8	1.78	0.49
37:3:113:C:C4	37:3:114:U:C4	3.00	0.49
37:3:91:G:H2'	37:3:92:A:H8	1.77	0.49
36:5:1253:U:O2	36:5:1263:A:H5'	2.12	0.49
36:5:1831:U:H2'	36:5:1832:C:H6	1.77	0.49
36:5:2147:A:H2'	36:5:2148:U:O4'	2.11	0.49
36:5:2881:C:H2'	36:5:2882:U:C6	2.48	0.49
1:6:406:U:H2'	1:6:407:A:C8	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:C1:33:ARG:NH1	13:C1:53:TYR:O	3.69	0.49
17:C5:98:ASN:HB2	17:C5:122:THR:HG22	1.95	0.49
1:2:1551:U:H3'	17:C5:43:ARG:NH2	2.27	0.49
19:C7:110:VAL:O	19:C7:114:GLY:N	2.46	0.49
28:D6:87:ARG:HD3	28:D6:92:ARG:HA	1.94	0.49
30:D8:8:THR:HB	30:D8:56:LEU:HB2	1.94	0.49
40:L3:283:TYR:HB3	40:L3:323:MET:CE	3.15	0.49
41:L4:334:PHE:HA	41:L4:339:LEU:HD11	2.79	0.49
42:L5:58:LYS:HA	42:L5:93:THR:HB	1.94	0.49
45:L8:70:LYS:HA	45:L8:235:GLY:HA3	3.82	0.49
46:L9:79:ILE:O	46:L9:82:VAL:HG12	2.38	0.49
52:M6:62:THR:HG21	52:M6:68:ARG:HG3	1.94	0.49
55:M9:168:ALA:C	55:M9:170:ARG:H	2.16	0.49
7:S5:172:ILE:O	7:S5:176:THR:HG23	2.58	0.49
36:1:2808:A:N7	36:1:2955:U:H4'	2.26	0.49
36:1:2864:A:O3'	47:M0:115:MET:HB2	2.11	0.49
1:2:641:G:H1	1:2:693:U:H3	1.61	0.49
36:5:1073:U:H2'	36:5:1074:U:C6	2.47	0.49
36:5:1194:G:O2'	36:5:1319:G:O2'	2.27	0.49
36:5:142:C:H2'	36:5:143:G:O4'	2.12	0.49
36:5:1472:U:H2'	36:5:1473:G:C8	2.47	0.49
36:5:2101:C:O2'	36:5:2102:U:OP1	2.31	0.49
36:5:2347:U:H2'	36:5:2348:A:O4'	2.13	0.49
36:5:2689:A:H2'	36:5:2689:A:N3	2.27	0.49
51:M5:172:ARG:NH1	36:5:29:C:O3'	106.06	0.49
36:5:3321:C:H2'	36:5:3322:A:C8	2.48	0.49
36:5:508:U:H2'	36:5:509:U:C6	2.47	0.49
1:6:1417:A:H2'	1:6:1418:G:O4'	2.12	0.49
1:6:5:U:H2'	1:6:6:G:H8	1.77	0.49
1:2:1389:C:O2'	19:C7:52:GLY:HA3	2.11	0.49
24:D2:23:ARG:NH1	24:D2:66:ASN:O	2.45	0.49
40:L3:4:ARG:CD	40:L3:7:GLU:HA	2.41	0.49
46:L9:48:VAL:HG11	46:L9:52:LEU:HD13	1.95	0.49
51:M5:28:TRP:CD1	36:5:2515:A:H5''	159.69	0.49
52:M6:32:LYS:O	52:M6:33:ILE:HD12	2.11	0.49
36:1:3309:G:H1'	53:M7:69:ARG:HD3	1.94	0.49
55:M9:158:GLU:HA	55:M9:161:ALA:HB3	2.70	0.49
59:N3:13:ILE:HG23	59:N3:85:TRP:CD1	2.47	0.49
77:Q1:13:LEU:HD11	77:Q1:17:ARG:CZ	2.42	0.49
8:S6:97:VAL:HG22	8:S6:98:ARG:H	2.28	0.49
19:C7:33:ARG:NH2	34:SR:109:ASP:OD1	2.97	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1027:A:H2'	36:1:1029:G:H5''	1.94	0.49
36:1:1103:A:H2'	36:1:1103:A:N3	2.28	0.49
36:1:2534:G:HO2'	36:1:2535:A:H8	1.60	0.49
36:1:2565:U:H2'	36:1:2566:C:H6	1.77	0.49
36:1:3217:C:C5	36:1:3220:G:H1'	2.47	0.49
1:2:709:C:N4	1:2:730:G:O2'	2.45	0.49
38:4:127:U:H2'	38:4:128:U:H5'	1.93	0.49
38:4:70:G:O6	73:O7:88:ALA:HB1	2.12	0.49
36:5:1946:A:H2'	36:5:1947:G:C8	2.47	0.49
51:M5:12:ARG:HG3	36:5:268:A:C4	127.37	0.49
36:5:284:A:H4'	36:5:285:A:C2	2.47	0.49
40:L3:334:ARG:NH2	36:5:3305:A:OP1	214.46	0.49
27:D5:77:ARG:HD2	1:6:1532:U:OP2	356.80	0.49
38:8:47:C:H1'	38:8:61:A:C4	2.47	0.49
38:8:82:U:O2	38:8:87:G:H4'	2.12	0.49
13:C1:35:TYR:HA	13:C1:61:THR:HA	2.72	0.49
13:C1:72:THR:H	13:C1:88:ARG:HH11	2.74	0.49
17:C5:56:PHE:O	17:C5:60:LEU:HB2	2.12	0.49
21:C9:115:GLU:O	21:C9:117:SER:N	2.45	0.49
40:L3:79:VAL:HG13	40:L3:322:ILE:HB	1.95	0.49
43:L6:82:ARG:O	69:O3:104:PRO:HA	2.86	0.49
46:L9:84:LYS:HZ3	46:L9:191:LEU:HD22	1.75	0.49
47:M0:189:GLU:HG3	47:M0:200:LEU:HD23	1.94	0.49
54:M8:150:VAL:HA	54:M8:153:PHE:CD1	2.47	0.49
54:M8:83:VAL:C	54:M8:85:GLY:H	2.41	0.49
57:N1:80:VAL:HG22	57:N1:83:ARG:HH21	2.05	0.49
36:1:1476:G:O3'	67:O1:63:GLY:HA2	2.12	0.49
70:O4:51:LEU:HD23	70:O4:51:LEU:H	1.77	0.49
3:S1:129:THR:HG22	3:S1:176:VAL:HG12	2.40	0.49
4:S2:139:ILE:HD12	4:S2:191:ALA:HB1	2.11	0.49
5:S3:75:LYS:HD2	12:C0:20:VAL:HG12	1.94	0.49
6:S4:9:LEU:HD12	6:S4:30:ARG:HA	1.95	0.49
7:S5:149:VAL:HG13	7:S5:151:GLY:H	5.53	0.49
7:S5:187:ILE:HD11	1:6:1535:U:H5''	330.39	0.49
7:S5:65:ARG:HE	7:S5:65:ARG:HA	4.55	0.49
34:SR:38:ARG:HG2	34:SR:67:ILE:HG23	1.95	0.49
36:1:1038:C:H4'	42:L5:5:LYS:HE3	1.94	0.49
36:1:408:A:N6	38:4:15:G:H1'	2.28	0.49
1:2:1376:C:N4	1:2:1377:U:O4	2.46	0.49
1:2:1474:G:H2'	1:2:1475:A:C8	2.48	0.49
1:2:1528:U:H2'	1:2:1529:C:H6	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:N1:130:ARG:HD3	36:5:1098:A:OP2	254.34	0.49
36:5:1932:A:H5'	36:5:1933:A:OP2	2.13	0.49
62:N6:12:ARG:HD3	36:5:215:G:H5''	87.63	0.49
36:5:2192:C:H2'	36:5:2193:U:O4'	2.12	0.49
36:5:2335:G:N2	36:5:2339:C:O2	2.38	0.49
36:5:394:G:N1	36:5:397:A:OP2	2.45	0.49
1:6:1243:G:N3	1:6:1243:G:H5''	2.27	0.49
1:6:1263:G:H2'	1:6:1264:G:O4'	2.13	0.49
1:6:1283:U:H2'	1:6:1284:C:C5	2.47	0.49
1:6:60:U:H5'	1:6:455:C:N4	2.28	0.49
1:6:607:G:H5'	1:6:613:G:N2	2.27	0.49
1:6:894:U:H2'	1:6:895:G:C8	2.48	0.49
20:C8:99:HIS:O	20:C8:101:LEU:HG	2.12	0.49
23:D1:53:TYR:CE2	23:D1:73:ALA:HB2	2.48	0.49
26:D4:51:GLU:O	26:D4:53:ASP:N	3.35	0.49
40:L3:14:LEU:HD22	40:L3:17:LEU:HD21	2.50	0.49
42:L5:40:HIS:CD2	42:L5:42:ALA:H	2.21	0.49
44:L7:132:PRO:HG2	44:L7:133:TYR:CE2	3.85	0.49
45:L8:48:ARG:HH21	45:L8:49:TYR:HE2	1.60	0.49
47:M0:52:LEU:HG	47:M0:152:LEU:HB3	1.95	0.49
54:M8:60:PRO:HG2	54:M8:142:GLY:HA3	3.45	0.49
54:M8:153:PHE:O	54:M8:161:LYS:HD3	2.13	0.49
62:N6:37:LYS:HA	62:N6:40:ARG:HB3	4.09	0.49
62:N6:82:VAL:O	62:N6:84:LYS:N	3.85	0.49
63:N7:14:VAL:HG13	70:O4:86:LYS:HG3	1.94	0.49
63:N7:14:VAL:N	63:N7:79:HIS:O	2.41	0.49
4:S2:176:SER:N	4:S2:195:ASP:OD1	2.46	0.49
1:2:66:U:C5	8:S6:173:PRO:HG3	2.48	0.49
1:2:767:U:C5	11:S9:143:ILE:HD12	2.48	0.49
11:S9:175:ARG:NH2	11:S9:179:ARG:HH22	2.10	0.49
36:1:1203:A:N6	36:1:1300:G:H2'	2.28	0.49
36:1:2152:A:HO2'	36:1:2243:A:HO2'	1.61	0.49
36:1:2947:G:OP2	40:L3:244:ARG:HD2	2.13	0.49
36:1:685:G:P	49:M3:35:ARG:HH11	2.36	0.49
1:2:1147:A:H2'	1:2:1148:C:H6	1.77	0.49
1:2:12:U:H2'	1:2:13:C:C6	2.48	0.49
1:2:1755:A:H2	1:2:1756[A]:A:C4	2.31	0.49
1:2:340:U:H2'	1:2:341:A:C8	2.47	0.49
1:2:615:A:O2'	1:2:621:A:N1	2.43	0.49
1:2:839:U:H2'	1:2:840:U:H5'	1.95	0.49
1:2:995:A:H2'	1:2:996:U:O4'	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:818:C:O2'	36:5:2138:A:N1	2.32	0.49
36:5:3159:C:H2'	36:5:3160:U:C6	2.48	0.49
1:6:1535:U:O2'	1:6:1536:G:O5'	2.30	0.49
1:6:16:G:H2'	1:6:17:C:C6	2.48	0.49
15:C3:54:LEU:HB3	15:C3:60:VAL:HG13	4.79	0.49
1:2:927:C:O2'	16:C4:125:SER:HB2	2.12	0.49
20:C8:139:LYS:O	20:C8:143:ARG:NH1	2.45	0.49
23:D1:12:TYR:CE2	23:D1:14:PRO:HG3	2.48	0.49
15:C3:61:THR:HG22	29:D7:32:PHE:CZ	3.59	0.49
41:L4:189:ALA:HA	36:5:1420:C:C5	116.06	0.49
42:L5:115:LEU:HD12	42:L5:119:TYR:HD2	3.84	0.49
44:L7:71:ALA:C	44:L7:73:GLY:H	2.14	0.49
45:L8:161:GLU:HA	45:L8:164:VAL:HG22	2.15	0.49
46:L9:120:ASP:OD1	46:L9:124:ARG:NH2	2.45	0.49
46:L9:41:ILE:HD11	46:L9:67:ALA:HB1	1.94	0.49
36:1:2853:A:O3'	47:M0:64:ALA:HB2	2.13	0.49
48:M1:53:THR:HG23	48:M1:59:ILE:O	2.12	0.49
36:1:3229:G:H1'	50:M4:133:LYS:HG2	1.93	0.49
52:M6:173:ALA:O	52:M6:176:LYS:HB3	2.75	0.49
63:N7:10:VAL:O	63:N7:83:THR:HG22	2.12	0.49
66:O0:103:THR:HG22	66:O0:104:LEU:H	4.09	0.49
71:O5:76:GLN:O	71:O5:81:ARG:NH1	2.45	0.49
72:O6:57:LEU:HD11	72:O6:73:ALA:HB2	1.97	0.49
72:O6:76:ARG:HA	72:O6:76:ARG:HE	2.02	0.49
5:S3:114:ALA:O	5:S3:116:ARG:N	3.14	0.49
34:SR:220:ILE:HB	34:SR:234:LEU:HB2	2.74	0.49
36:1:2900:A:N3	36:1:3025:C:O2'	2.37	0.49
36:1:304:G:C6	51:M5:178:HIS:CD2	3.01	0.49
36:1:577:C:O2'	36:1:579:G:OP1	2.30	0.49
36:1:656:A:H2'	36:1:657:A:C8	2.48	0.49
1:2:1254:U:OP2	14:C2:46:ARG:NH2	2.44	0.49
1:2:1415:U:OP1	19:C7:45:ARG:NH2	2.40	0.49
1:2:802:G:H21	24:D2:107:SER:HB3	1.78	0.49
36:5:1577:G:H2'	36:5:1578:C:C6	2.48	0.49
36:5:2599:U:H2'	36:5:2600:C:C6	2.48	0.49
36:5:279:U:H2'	36:5:280:U:H6	1.78	0.49
36:5:3242:G:N2	36:5:3245:A:H5''	2.28	0.49
36:5:3279:A:N6	36:5:3280:U:O4	2.46	0.49
36:5:3359:A:H2'	36:5:3360:C:O4'	2.12	0.49
36:5:438:A:H3'	36:5:439:C:C6	2.47	0.49
36:5:627:U:H2'	36:5:628:A:C8	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1394:G:H2'	1:6:1395:G:C8	2.48	0.49
1:6:196:G:N3	1:6:197:A:H1'	2.28	0.49
14:C2:128:ALA:HB3	14:C2:133:LEU:HD22	3.50	0.49
15:C3:66:ILE:HG13	15:C3:67:THR:HG23	1.95	0.49
2:S0:66:ALA:HB2	23:D1:37:ALA:HB2	2.27	0.49
24:D2:81:VAL:HG11	24:D2:86:ILE:HG23	1.95	0.49
24:D2:9:ASP:HA	24:D2:12:ASN:HB3	2.87	0.49
40:L3:173:GLN:O	40:L3:173:GLN:HG3	2.13	0.49
43:L6:40:LEU:HB3	43:L6:84:VAL:HG11	3.93	0.49
46:L9:128:VAL:HG13	46:L9:134:ILE:HD12	1.94	0.49
47:M0:9:TYR:CG	47:M0:97:LEU:HD13	2.52	0.49
49:M3:76:THR:HG21	49:M3:103:ASN:OD1	3.34	0.49
49:M3:113:VAL:HG12	49:M3:117:LYS:HE3	3.17	0.49
55:M9:125:LYS:NZ	36:5:1720:U:O4	240.81	0.49
56:N0:12:ARG:NH1	56:N0:15:PRO:HG3	2.28	0.49
59:N3:32:ARG:HB3	59:N3:64:LYS:O	2.12	0.49
59:N3:45:ARG:HD2	59:N3:46:LEU:H	2.38	0.49
61:N5:41:ALA:O	61:N5:43:ALA:N	3.63	0.49
3:S1:35:PRO:HD3	3:S1:98:THR:HG23	1.94	0.49
4:S2:148:LEU:HD13	4:S2:149:GLY:H	1.77	0.49
4:S2:44:LEU:HD21	4:S2:247:ALA:HB2	2.14	0.49
9:S7:58:LEU:HG	9:S7:88:ARG:HD2	1.94	0.49
10:S8:25:ARG:O	10:S8:28:GLU:HG2	2.15	0.49
11:S9:100:LYS:HE3	11:S9:103:ASP:OD2	2.12	0.49
34:SR:47:LEU:HD22	34:SR:54:PHE:CE2	4.28	0.49
36:1:1471:U:H2'	36:1:1472:U:C6	2.47	0.49
36:1:2359:C:H2'	36:1:2360:C:C6	2.47	0.49
36:1:2651:G:H4'	36:1:2652:U:OP2	2.13	0.49
1:2:972:G:O2'	36:1:847:A:N1	2.32	0.49
1:2:1488:G:N3	1:2:1495:C:H1'	2.28	0.49
1:2:386:G:OP2	10:S8:25:ARG:NH2	2.45	0.49
1:2:877:G:H5'	1:2:937:C:H1'	1.95	0.49
36:5:247:C:N4	36:5:248:U:O2	2.45	0.49
36:5:3212:C:H2'	36:5:3213:A:O4'	2.12	0.49
36:5:750:G:H2'	36:5:751:A:H8	1.77	0.49
1:6:108:A:H2'	1:6:109:G:C8	2.47	0.49
1:6:1175:U:H2'	1:6:1176:G:H8	1.76	0.49
1:6:1628:U:H2'	1:6:1629:G:C8	2.48	0.49
1:6:1645:G:H1	1:6:1756[A]:A:H61	1.61	0.49
1:6:721:U:O2'	1:6:722:G:O4'	2.29	0.49
40:L3:292:ALA:HA	40:L3:303:LYS:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:L3:280:HIS:HB3	40:L3:324:VAL:CG1	2.42	0.49
43:L6:28:GLN:HG3	36:5:501:A:O3'	257.70	0.49
36:1:1334:U:H5''	44:L7:206:LYS:HB3	1.95	0.49
48:M1:143:ARG:HG2	48:M1:144:CYS:SG	2.53	0.49
50:M4:47:ASP:HB2	50:M4:55:ARG:HG3	2.66	0.49
53:M7:138:LYS:HG3	53:M7:140:GLU:HG3	1.94	0.49
36:1:840:C:H4'	55:M9:128:LYS:HD3	1.94	0.49
56:N0:65:ASN:ND2	36:5:521:A:N3	315.04	0.49
57:N1:79:MET:HA	57:N1:84:TYR:HA	1.93	0.49
49:M3:155:GLU:OE2	64:N8:90:TYR:OH	4.81	0.49
69:O3:37:THR:HG23	69:O3:40:ASP:HB2	1.95	0.49
70:O4:107:GLU:HA	70:O4:110:GLU:CD	4.55	0.49
72:O6:90:MET:SD	72:O6:90:MET:N	2.86	0.49
2:S0:53:THR:HA	2:S0:161:PRO:HG2	1.95	0.49
6:S4:33:ALA:O	1:6:121:U:O2'	352.55	0.49
10:S8:137:LYS:HG2	10:S8:138:ASN:H	1.77	0.49
10:S8:103:GLN:HA	10:S8:165:LEU:O	2.61	0.49
10:S8:183:ILE:HG13	10:S8:183:ILE:O	4.73	0.49
11:S9:54:ARG:HH11	11:S9:55:ALA:HB2	1.78	0.49
34:SR:273:ASP:OD2	34:SR:275:ARG:NH1	4.48	0.49
36:1:1355:A:H5''	36:1:1356:U:C5	2.44	0.49
36:1:1723:A:OP1	55:M9:128:LYS:NZ	2.44	0.49
36:1:2180:G:H2'	36:1:2181:C:C6	2.48	0.49
36:1:2726:C:O2'	36:1:2727:A:H2'	2.13	0.49
36:1:729:C:H2'	36:1:730:C:C6	2.48	0.49
1:2:1277:G:H2'	1:2:1278:G:O4'	2.12	0.49
1:2:292:U:H2'	1:2:293:U:H6	1.78	0.49
36:5:1239:C:N4	36:5:1249:G:H22	2.11	0.49
36:5:3160:U:H3	36:5:3290:G:H1	1.60	0.49
33:E1:135:HIS:HB3	1:6:1250:U:H2'	431.95	0.49
19:C7:45:ARG:HD3	1:6:1332:C:OP2	420.31	0.49
1:6:271:A:C2	1:6:285:G:C6	3.01	0.49
1:6:845:G:H2'	1:6:846:G:C8	2.41	0.49
13:C1:57:LYS:HB2	13:C1:110:HIS:NE2	2.27	0.49
20:C8:28:ILE:HG23	20:C8:58:ALA:HA	5.16	0.49
20:C8:56:LYS:HD3	20:C8:61:LEU:HD23	1.94	0.49
24:D2:9:ASP:HB3	1:6:1036:A:O2'	361.70	0.49
1:2:1433:G:N1	31:D9:45:GLU:OE2	2.46	0.49
33:E1:82:LYS:C	33:E1:84:VAL:H	4.28	0.49
39:L2:80:GLU:N	39:L2:170:ALA:HB2	3.67	0.49
40:L3:103:THR:HG21	40:L3:147:GLU:OE2	2.22	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:L4:6:VAL:O	41:L4:20:LEU:N	2.45	0.49
45:L8:137:ASN:ND2	36:5:148:G:N7	108.79	0.49
45:L8:27:THR:HG22	63:N7:53:VAL:HG12	7.13	0.49
47:M0:44:ASP:CG	47:M0:185:ARG:HH11	2.16	0.49
48:M1:20:ASN:HB3	48:M1:126:ASP:HB2	1.95	0.49
52:M6:9:ILE:O	52:M6:36:VAL:HG22	2.13	0.49
60:N4:97:LYS:H	60:N4:98:PRO:HD2	1.78	0.49
62:N6:57:LEU:HB3	62:N6:105:VAL:HG12	2.80	0.49
64:N8:46:ASP:N	64:N8:46:ASP:OD1	2.45	0.49
36:1:3383:G:N2	67:O1:105:GLN:OE1	2.33	0.49
36:1:424:G:O2'	68:O2:23:ASP:OD2	2.28	0.49
71:O5:107:LYS:HA	71:O5:110:ALA:HB3	3.10	0.49
8:S6:135:PRO:HB2	8:S6:141:ILE:HD13	4.14	0.49
11:S9:114:TYR:HA	11:S9:119:ALA:HB3	2.22	0.49
11:S9:166:GLY:O	11:S9:168:ARG:N	2.46	0.49
35:SM:35:ALA:HB1	35:SM:37:VAL:HG23	1.94	0.49
34:SR:243:LEU:HD22	34:SR:252:LEU:HD11	1.95	0.49
34:SR:69:GLN:OE1	34:SR:85:TRP:NE1	2.40	0.49
36:1:649:A:OP2	36:1:2868:U:O2'	2.22	0.48
1:2:1256:A:OP1	12:C0:5:LYS:NZ	2.29	0.48
1:2:1655:A:N1	36:1:2291:A:O2'	2.39	0.48
1:2:711:U:H4'	1:2:712:G:OP1	2.13	0.48
46:L9:63:LYS:NZ	36:5:1210:U:H5'	312.83	0.48
36:5:1260:A:O2'	36:5:1279:C:O2	2.31	0.48
36:5:1497:C:H2'	36:5:1498:A:C8	2.48	0.48
39:L2:241:ARG:HG2	36:5:2155:G:OP1	221.69	0.48
72:O6:28:TYR:HA	36:5:316:U:O4	101.75	0.48
1:6:1370:U:H4'	1:6:1371:A:H4'	1.94	0.48
1:6:1624:C:H2'	1:6:1625:C:C6	2.48	0.48
1:6:1642:G:H2'	1:6:1643:U:H6	1.78	0.48
1:6:417:A:H4'	1:6:418:G:O5'	2.13	0.48
13:C1:72:THR:H	13:C1:88:ARG:NH1	3.41	0.48
13:C1:5:LEU:HG	13:C1:9:SER:OG	8.85	0.48
16:C4:29:HIS:ND1	16:C4:29:HIS:O	2.46	0.48
20:C8:126:ARG:HG2	20:C8:133:VAL:HA	1.95	0.48
28:D6:89:ARG:HD2	28:D6:92:ARG:NH2	2.28	0.48
40:L3:332:ARG:O	40:L3:333:LYS:HB2	2.28	0.48
41:L4:311:HIS:ND1	44:L7:162:PRO:HG2	2.28	0.48
44:L7:90:LYS:HD2	44:L7:91:GLY:H	1.77	0.48
47:M0:76:MET:HE1	47:M0:148:VAL:HG22	2.78	0.48
47:M0:99:ILE:HD12	47:M0:123:HIS:ND1	3.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:M1:139:THR:HG22	48:M1:147:THR:HA	1.95	0.48
53:M7:52:LEU:HD11	53:M7:88:VAL:HB	1.95	0.48
57:N1:39:ILE:HG13	57:N1:102:ARG:HD2	5.34	0.48
60:N4:47:ARG:HD3	60:N4:58:HIS:CD2	2.48	0.48
63:N7:10:VAL:HG23	63:N7:86:THR:HA	1.95	0.48
36:1:1407:A:O3'	68:O2:33:ARG:NH2	2.46	0.48
79:Q3:86:LEU:O	79:Q3:90:VAL:HG13	6.55	0.48
6:S4:38:LEU:HB3	1:6:298:C:H5''	352.10	0.48
36:1:2107:A:H2	36:1:3344:A:C8	2.31	0.48
36:1:776:U:H5	36:1:2719:U:O2	1.97	0.48
36:1:3244:A:OP1	40:L3:97:ARG:NH2	2.46	0.48
36:1:3274:A:H2'	53:M7:171:ARG:HH12	1.79	0.48
1:2:1142:A:H2'	1:2:1143:A:C8	2.48	0.48
1:2:1202:A:H1'	1:2:1207:C:H42	1.77	0.48
1:2:1573:A:H4'	1:2:1574:G:OP2	2.12	0.48
1:2:1719:A:H2'	1:2:1720:G:O4'	2.13	0.48
1:2:6:G:H2'	1:2:7:G:H8	1.78	0.48
36:5:2584:G:H5'	36:5:2585:G:OP2	2.14	0.48
36:5:2818:U:H6	36:5:2818:U:H5'	1.77	0.48
36:5:3006:A:H2'	36:5:3007:U:O4'	2.12	0.48
36:5:3060:C:C2	36:5:3061:G:C8	3.01	0.48
36:5:406:G:H1'	38:8:16:G:N2	2.29	0.48
69:O3:45:LEU:HD12	36:5:584:G:O2'	253.15	0.48
36:5:638:C:H2'	36:5:639:G:C8	2.48	0.48
1:6:691:C:OP1	1:6:696:C:N4	2.39	0.48
15:C3:87:ASP:OD1	1:6:867:G:N2	317.37	0.48
29:D7:28:PRO:HB3	1:6:959:U:H5'	350.83	0.48
38:8:121:U:H2'	38:8:122:U:C6	2.48	0.48
18:C6:40:GLU:HG3	18:C6:42:GLU:HB3	1.95	0.48
36:1:2620:G:H22	84:C:74:CH:HN3	1.59	0.48
22:D0:104:THR:HG22	22:D0:116:VAL:HG21	1.95	0.48
25:D3:16:ARG:O	25:D3:19:ARG:HB3	2.13	0.48
26:D4:68:LYS:HE2	26:D4:70:VAL:HG22	3.50	0.48
41:L4:193:LYS:O	41:L4:198:ARG:HG2	3.78	0.48
41:L4:271:LYS:HB2	41:L4:274:TYR:HB3	2.22	0.48
42:L5:119:TYR:OH	42:L5:139:PRO:O	2.82	0.48
45:L8:182:GLY:O	45:L8:186:LEU:HG	2.39	0.48
50:M4:94:TRP:CE2	50:M4:100:ALA:HB2	2.48	0.48
51:M5:11:GLN:HG2	51:M5:44:ARG:HH21	2.23	0.48
36:1:3243:A:C8	52:M6:156:LEU:HD22	2.48	0.48
54:M8:154:GLY:O	54:M8:159:LYS:HE3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:M9:161:ALA:O	55:M9:165:LYS:HG2	2.12	0.48
66:O0:66:LYS:N	66:O0:66:LYS:HD2	4.39	0.48
69:O3:14:LEU:HD11	69:O3:31:LYS:HB2	2.69	0.48
78:Q2:41:ARG:HH12	36:5:284:A:P	155.73	0.48
79:Q3:13:LYS:HD2	79:Q3:14:TYR:CZ	3.23	0.48
4:S2:144:TRP:CE2	4:S2:173:PRO:HG3	3.11	0.48
6:S4:163:ASP:O	6:S4:165:ALA:N	2.44	0.48
6:S4:88:ASP:OD1	6:S4:122:LYS:NZ	2.38	0.48
7:S5:45:LYS:HD3	7:S5:45:LYS:HA	1.81	0.48
9:S7:162:ILE:HA	9:S7:165:LYS:HD2	1.93	0.48
9:S7:165:LYS:O	9:S7:168:SER:OG	2.27	0.48
10:S8:33:PRO:HB3	1:6:330:G:O2'	273.36	0.48
36:1:2097:U:H2'	36:1:2098:C:C6	2.48	0.48
36:1:543:C:H3'	36:1:544:C:C6	2.48	0.48
1:2:1250:U:HO2'	1:2:1251:U:P	2.36	0.48
1:2:851:U:H2'	1:2:852:C:C6	2.48	0.48
36:5:1312:C:H2'	36:5:1313:G:O4'	2.12	0.48
36:5:1658:G:H2'	36:5:1659:U:C6	2.48	0.48
36:5:1794:G:O2'	36:5:1795:U:H5'	2.14	0.48
70:O4:63:ALA:HB2	36:5:1803:C:H5'	157.97	0.48
36:5:1131:G:C4	36:5:2373:A:C2	3.01	0.48
36:5:646:A:C2	36:5:2375:G:C2	3.01	0.48
64:N8:36:GLY:N	36:5:40:A:OP2	174.01	0.48
1:6:1427:A:O2'	1:6:1428:G:OP1	2.30	0.48
27:D5:96:SER:OG	1:6:1530:C:OP2	371.64	0.48
1:6:712:G:H2'	1:6:713:A:H8	1.79	0.48
12:C0:29:GLN:O	12:C0:31:LYS:N	2.43	0.48
14:C2:46:ARG:HG3	33:E1:103:LEU:HD12	1.94	0.48
1:2:867:G:OP2	15:C3:3:ARG:NH1	2.46	0.48
15:C3:41:ALA:HB1	15:C3:75:LEU:HD21	3.39	0.48
17:C5:130:ARG:N	17:C5:130:ARG:HH11	2.07	0.48
19:C7:58:MET:O	19:C7:61:ILE:HG22	5.19	0.48
21:C9:108:LEU:HB3	21:C9:114:VAL:HG22	6.39	0.48
21:C9:141:GLU:C	21:C9:143:ASP:H	3.26	0.48
1:2:1479:A:H5''	21:C9:60:SER:OG	2.13	0.48
21:C9:61:VAL:O	21:C9:65:ILE:HG13	2.29	0.48
23:D1:3:ASN:ND2	23:D1:6:GLY:O	2.45	0.48
25:D3:109:ARG:NH2	25:D3:113:ALA:O	2.46	0.48
1:2:1199:G:C5	31:D9:40:ARG:HD3	2.48	0.48
40:L3:250:ALA:HB3	36:5:2880:U:O2	223.81	0.48
40:L3:221:THR:HB	40:L3:273:HIS:H	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:L6:100:LYS:NZ	43:L6:137:ASP:OD2	2.41	0.48
47:M0:3:ARG:NH2	36:5:2854:U:OP2	290.44	0.48
45:L8:136:LEU:HD22	51:M5:3:ALA:HB2	2.18	0.48
61:N5:88:MET:HA	61:N5:120:LYS:HG3	1.95	0.48
62:N6:57:LEU:HD22	62:N6:58:VAL:H	2.93	0.48
63:N7:11:ALA:HB1	63:N7:80:LEU:HB2	1.94	0.48
36:1:1117:G:OP1	65:N9:4:SER:HB2	2.14	0.48
66:O0:99:ASP:N	66:O0:99:ASP:OD2	2.45	0.48
70:O4:102:LYS:O	70:O4:106:LYS:HD3	5.82	0.48
70:O4:83:ASN:OD1	70:O4:83:ASN:N	3.29	0.48
71:O5:17:LEU:HA	71:O5:20:GLN:HB2	2.69	0.48
72:O6:62:ARG:HH21	72:O6:94:ILE:HG13	6.74	0.48
79:Q3:36:ARG:O	79:Q3:45:LYS:NZ	2.37	0.48
7:S5:66:GLN:HG3	7:S5:67:PRO:HD2	1.94	0.48
36:1:1430:U:O4	64:N8:3:SER:OG	2.31	0.48
36:1:1767:C:H2'	36:1:1768:U:C6	2.48	0.48
36:1:1832:C:H2'	36:1:1833:G:H8	1.79	0.48
36:1:660:A:H5''	41:L4:100:PHE:CD1	2.47	0.48
1:2:1041:G:H2'	1:2:1042:G:C8	2.48	0.48
1:2:886:U:H2'	1:2:887:A:O4'	2.14	0.48
36:5:1313:G:H2'	36:5:1314:C:H6	1.78	0.48
70:O4:41:ARG:HD3	36:5:1739:U:H1'	187.38	0.48
36:5:2370:G:H2'	36:5:2371:G:O4'	2.13	0.48
45:L8:46:LEU:HD21	36:5:2524:A:C4	182.33	0.48
36:5:2526:C:H1'	36:5:2588:U:H5''	1.94	0.48
36:5:3242:G:N2	36:5:3245:A:OP2	2.44	0.48
36:5:3306:U:O2'	36:5:3308:C:OP2	2.21	0.48
36:5:3378:C:H2'	36:5:3379:C:C6	2.48	0.48
54:M8:142:GLY:O	36:5:744:A:H4'	168.09	0.48
5:S3:162:GLN:HG3	1:6:1333:C:H4'	427.39	0.48
1:6:414:C:H2'	1:6:415:C:O4'	2.14	0.48
1:6:855:A:C2	1:6:857:U:H1'	2.48	0.48
12:C0:30:ALA:O	12:C0:38:LYS:HA	2.13	0.48
21:C9:105:LEU:HB3	21:C9:122:ARG:NE	2.29	0.48
1:2:522:U:H5''	26:D4:37:LYS:HG3	1.94	0.48
40:L3:169:THR:HG22	40:L3:171:LEU:HG	1.94	0.48
41:L4:262:TRP:CZ3	41:L4:271:LYS:HE3	2.93	0.48
36:1:515:C:O2'	41:L4:342:LYS:O	2.31	0.48
46:L9:19:SER:HA	50:M4:6:ILE:O	2.48	0.48
47:M0:47:PRO:HB3	47:M0:171:TRP:CE2	2.71	0.48
56:N0:77:VAL:N	56:N0:92:LYS:O	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
58:N2:67:SER:OG	58:N2:68:THR:N	2.46	0.48
60:N4:49:ILE:O	60:N4:52:THR:OG1	2.21	0.48
36:1:359:U:HO2'	73:O7:16:HIS:HD1	1.58	0.48
75:O9:27:ILE:HG23	75:O9:30:ARG:CZ	3.86	0.48
76:Q0:106:ARG:NH1	76:Q0:106:ARG:HB2	4.61	0.48
3:S1:222:LYS:HG3	3:S1:223:PHE:H	1.78	0.48
4:S2:88:LYS:HD2	4:S2:95:ARG:CZ	8.73	0.48
6:S4:100:ARG:NH2	6:S4:121:TYR:O	2.46	0.48
36:1:1752:A:H8	36:1:1752:A:O5'	1.96	0.48
36:1:3133:C:H2'	36:1:3134:A:O4'	2.14	0.48
36:1:345:G:OP1	36:1:1429:G:N1	2.37	0.48
36:1:644:G:H2'	36:1:2372:A:N7	2.28	0.48
36:1:896:A:H5'	39:L2:183:GLY:HA2	1.95	0.48
38:4:58:G:O6	73:O7:63:ARG:NH2	2.40	0.48
36:5:1070:U:H2'	36:5:1071:U:O4'	2.12	0.48
36:5:1759:C:H2'	36:5:1760:A:O4'	2.14	0.48
52:M6:68:ARG:NH1	36:5:2988:C:P	216.62	0.48
55:M9:59:SER:N	36:5:3068:U:OP1	165.36	0.48
1:6:498:G:N7	1:6:499:U:N3	2.62	0.48
1:6:716:C:H1'	1:6:723:G:H22	1.79	0.48
1:6:955:A:H4'	1:6:1073:G:O2'	2.13	0.48
17:C5:78:THR:OG1	17:C5:79:HIS:N	2.86	0.48
1:2:1459:C:N4	20:C8:139:LYS:HE2	2.29	0.48
21:C9:125:SER:O	21:C9:129:GLN:HG3	2.13	0.48
23:D1:77:GLY:O	23:D1:78:LEU:HD12	6.54	0.48
25:D3:34:LEU:O	25:D3:39:LYS:NZ	2.47	0.48
39:L2:62:VAL:HG21	39:L2:71:LEU:HD23	1.95	0.48
40:L3:283:TYR:HB3	40:L3:356:LEU:HD21	1.96	0.48
41:L4:20:LEU:HD11	41:L4:252:GLU:HG3	3.78	0.48
44:L7:131:GLU:HG3	44:L7:230:GLY:HA2	4.68	0.48
37:3:43:U:H4'	48:M1:140:ARG:O	2.14	0.48
50:M4:132:LYS:HD3	36:5:3230:G:H4'	286.50	0.48
52:M6:35:VAL:HB	52:M6:104:VAL:HG13	1.96	0.48
52:M6:28:LEU:HD21	52:M6:88:VAL:HG13	1.93	0.48
53:M7:108:ASP:N	53:M7:152:GLU:OE2	2.62	0.48
61:N5:67:ILE:HD12	61:N5:83:VAL:HG12	2.51	0.48
70:O4:86:LYS:O	70:O4:90:ILE:HG12	2.13	0.48
9:S7:51:VAL:HG23	9:S7:53:GLY:N	2.28	0.48
11:S9:6:ARG:NH1	1:6:39:A:OP1	385.31	0.48
36:1:1134:G:C2	36:1:1135:A:C8	3.01	0.48
36:1:1703:U:N3	36:1:1740:U:O2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:2254:U:H2'	36:1:2261:G:H22	1.78	0.48
36:1:2597:U:H2'	36:1:2598:G:H8	1.77	0.48
36:1:2807:U:O2'	36:1:2809:C:OP1	2.30	0.48
36:1:563:U:H2'	36:1:564:G:H8	1.79	0.48
36:1:920:A:OP1	36:1:923:C:N4	2.43	0.48
1:2:1162:C:H1'	1:2:1616:G:N2	2.29	0.48
1:2:735:C:O2'	1:2:736:C:H5''	2.14	0.48
1:2:826:U:H2'	1:2:827:C:H6	1.78	0.48
38:4:24:G:OP2	62:N6:13:ARG:HD3	2.13	0.48
53:M7:67:ILE:HD11	36:5:1447:G:H3'	165.11	0.48
36:5:2372:A:H4'	36:5:2373:A:OP2	2.13	0.48
36:5:2403:G:N7	36:5:2870:C:H4'	2.28	0.48
36:5:2533:G:H21	36:5:2546:C:H42	1.61	0.48
36:5:2572:C:O2'	36:5:2573:G:H5''	2.13	0.48
36:5:3356:G:C6	36:5:3357:U:C4	3.01	0.48
36:5:3386:G:H2'	36:5:3387:U:C6	2.49	0.48
64:N8:33:GLY:N	36:5:798:G:OP1	158.72	0.48
1:6:250:C:H2'	1:6:251:A:C8	2.49	0.48
19:C7:26:LEU:HD13	19:C7:59:LYS:HG3	1.95	0.48
20:C8:83:ALA:HA	20:C8:86:LEU:HD22	1.94	0.48
26:D4:10:ARG:HG3	26:D4:26:ASP:OD2	4.88	0.48
26:D4:80:ALA:HA	26:D4:83:LYS:HE2	6.19	0.48
29:D7:34:ASP:HB3	29:D7:43:ILE:HD12	1.96	0.48
1:2:1553:G:H4'	31:D9:14:TYR:CE1	2.45	0.48
39:L2:183:GLY:O	39:L2:186:PHE:HB3	2.22	0.48
36:1:2941:A:OP2	40:L3:256:HIS:HD2	1.97	0.48
41:L4:173:GLY:O	41:L4:175:HIS:N	3.09	0.48
42:L5:211:LEU:HD22	42:L5:211:LEU:HA	1.72	0.48
37:3:49:G:N7	42:L5:58:LYS:HE2	2.28	0.48
44:L7:178:ILE:HA	44:L7:183:ASP:HB3	1.94	0.48
46:L9:86:TYR:CD2	46:L9:151:VAL:HG22	2.47	0.48
49:M3:76:THR:OG1	49:M3:77:LEU:N	3.72	0.48
55:M9:20:ARG:HG2	36:5:1875:G:OP2	136.95	0.48
55:M9:8:LYS:HG3	55:M9:22:VAL:HG21	1.96	0.48
36:1:662:U:OP1	64:N8:8:THR:HG21	2.12	0.48
67:O1:79:ARG:NE	67:O1:79:ARG:H	2.10	0.48
76:Q0:110:CYS:SG	76:Q0:112:LYS:HB2	2.96	0.48
78:Q2:43:TYR:OH	78:Q2:47:GLN:NE2	2.47	0.48
2:S0:180:GLU:O	2:S0:184:LEU:HD22	3.08	0.48
3:S1:51:SER:HA	3:S1:57:ALA:H	1.79	0.48
5:S3:34:TYR:OH	5:S3:37:VAL:HG13	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S3:94:ARG:HH21	35:SM:134:ASP:CG	2.07	0.48
6:S4:114:ILE:HD12	6:S4:118:GLU:HG2	2.66	0.48
1:2:446:A:H5'	6:S4:57:ASN:HB3	1.96	0.48
9:S7:49:ILE:HD11	9:S7:172:VAL:HG13	3.00	0.48
11:S9:40:LYS:HA	11:S9:43:TYR:HB2	2.70	0.48
34:SR:106:HIS:HD2	34:SR:110:VAL:HG22	1.96	0.48
36:1:1362:G:O2'	44:L7:159:GLN:HA	2.13	0.48
36:1:1734:G:H2'	36:1:1735:G:O4'	2.14	0.48
36:1:2911:A:H4'	36:1:2912:G:C8	2.47	0.48
36:1:379:C:H2'	36:1:380:U:C6	2.49	0.48
36:1:381:U:H2'	36:1:382:U:C6	2.49	0.48
36:1:437:G:H22	36:1:622:A:H61	1.62	0.48
36:1:685:G:OP1	49:M3:35:ARG:HD2	2.14	0.48
36:1:686:G:C6	36:1:687:U:C2	3.02	0.48
1:2:1146:G:C6	1:2:1147:A:C6	3.01	0.48
1:2:16:G:H2'	1:2:17:C:C6	2.47	0.48
36:5:1846:C:H3'	36:5:1847:A:H8	1.78	0.48
36:5:550:A:H2'	36:5:551:A:C8	2.48	0.48
41:L4:310:THR:HG22	36:5:609:G:C8	225.78	0.48
54:M8:43:PRO:HG2	36:5:729:C:P	191.64	0.48
6:S4:6:LYS:NZ	1:6:383:G:OP1	344.19	0.48
15:C3:139:TRP:HZ2	15:C3:149:LEU:HD21	1.78	0.48
1:2:1788:G:P	16:C4:127:ARG:HH22	2.36	0.48
1:2:1453:G:H5'	17:C5:79:HIS:HB3	1.95	0.48
18:C6:20:ALA:HB2	18:C6:67:VAL:HG13	1.96	0.48
26:D4:37:LYS:HG3	1:6:522:U:H5''	419.06	0.48
32:E0:49:LEU:H	32:E0:49:LEU:HD22	2.80	0.48
39:L2:140:ASN:HD21	39:L2:142:ASP:HB3	6.07	0.48
40:L3:116:ARG:NH2	40:L3:174:LYS:HD3	2.28	0.48
40:L3:148:LEU:HD21	40:L3:196:ARG:HB2	1.96	0.48
41:L4:60:THR:HG22	41:L4:62:ALA:N	2.76	0.48
42:L5:183:TRP:CZ3	42:L5:185:PHE:HA	6.34	0.48
43:L6:41:ILE:HB	43:L6:85:ILE:HB	2.09	0.48
45:L8:100:GLU:OE1	45:L8:108:ARG:HD3	2.98	0.48
46:L9:77:ASN:N	46:L9:77:ASN:OD1	2.46	0.48
57:N1:36:VAL:HG13	57:N1:65:TYR:HA	1.95	0.48
64:N8:91:LEU:HD13	64:N8:121:VAL:HG21	1.94	0.48
3:S1:27:LYS:NZ	3:S1:49:ASN:OD1	4.82	0.48
3:S1:68:VAL:HG23	3:S1:73:LEU:HD21	5.36	0.48
6:S4:43:PRO:HA	6:S4:82:TYR:O	2.46	0.48
9:S7:109:VAL:HG22	9:S7:110:GLN:H	4.23	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:S8:42:ARG:O	10:S8:58:LEU:HB2	4.87	0.48
11:S9:31:ALA:O	11:S9:36:LEU:N	3.19	0.48
36:1:1076:C:O3'	65:N9:38:LYS:NZ	2.43	0.48
36:1:1571:A:H2'	36:1:1572:U:O4'	2.14	0.48
36:1:1791:C:H2'	36:1:1792:C:C6	2.49	0.48
36:1:2553:U:C4	70:O4:95:ILE:HG12	2.49	0.48
36:1:2723:U:H2'	36:1:2724:U:C6	2.49	0.48
36:1:3013:U:H2'	36:1:3014:U:C6	2.49	0.48
36:1:3204:C:O2'	36:1:3205:G:H5'	2.13	0.48
1:2:1338:C:H1'	1:2:1410:A:C4	2.48	0.48
1:2:749:U:H2'	1:2:750:U:C6	2.49	0.48
1:2:993:A:H2'	1:2:994:G:O4'	2.14	0.48
36:5:1918:C:H2'	36:5:1919:G:O4'	2.14	0.48
36:5:2568:C:N3	36:5:2574:G:N1	2.61	0.48
36:5:2993:G:C6	36:5:3142:A:C4	3.01	0.48
1:6:249:U:H3'	1:6:250:C:H5'	1.94	0.48
32:E0:26:LYS:NZ	1:6:588:U:OP1	421.51	0.48
16:C4:123:SER:HB2	1:6:885:G:H21	285.62	0.48
22:D0:95:ALA:HB1	22:D0:99:ILE:HB	4.38	0.48
24:D2:37:PHE:CE1	24:D2:103:ILE:HD12	4.31	0.48
41:L4:300:ARG:HG2	54:M8:39:ARG:O	2.27	0.48
41:L4:316:ASN:HB3	41:L4:319:LYS:HG2	1.95	0.48
42:L5:34:LYS:HE2	42:L5:38:THR:OG1	7.43	0.48
47:M0:97:LEU:O	47:M0:123:HIS:N	2.57	0.48
52:M6:110:PRO:O	52:M6:112:TYR:N	2.77	0.48
53:M7:32:THR:HG22	53:M7:58:ILE:HG21	1.94	0.48
56:N0:27:MET:HE2	56:N0:29:ILE:HD11	1.96	0.48
65:N9:25:LYS:NZ	36:5:1107:C:H5''	198.49	0.48
68:O2:101:SER:O	68:O2:105:ARG:HG3	2.13	0.48
5:S3:105:MET:HG2	5:S3:122:VAL:HG21	4.58	0.48
6:S4:123:LEU:HD23	6:S4:228:ILE:HG22	1.94	0.48
9:S7:99:LEU:HD23	9:S7:116:ARG:HB3	4.88	0.48
9:S7:138:LYS:HE3	9:S7:138:LYS:HB2	1.67	0.48
10:S8:61:GLU:HG3	10:S8:62:THR:HG23	1.94	0.48
11:S9:62:ARG:NH1	11:S9:68:LYS:HD3	2.69	0.48
34:SR:31:ASN:HA	34:SR:47:LEU:HD12	3.75	0.48
36:1:1460:A:H2'	36:1:1461:A:C8	2.49	0.48
36:1:1525:G:H5'	36:1:1830:G:OP2	2.14	0.48
36:1:1711:C:H2'	36:1:1712:G:O4'	2.14	0.48
36:1:2144:A:H1'	36:1:2281:A:N6	2.28	0.48
36:1:76:G:O2'	49:M3:100:ARG:NH1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1250:U:O2'	1:2:1251:U:OP1	2.27	0.48
1:2:1396:U:H2'	1:2:1397:U:C6	2.48	0.48
1:2:158:U:HO2'	1:2:159:U:H3'	1.78	0.48
1:2:1657:U:H1'	1:2:1658:G:OP2	2.14	0.48
1:2:934:C:N3	1:2:1077:C:H4'	2.28	0.48
37:3:61:G:H2'	37:3:62:U:C6	2.49	0.48
36:5:122:A:H5''	36:5:123:A:C8	2.49	0.48
45:L8:136:LEU:HD12	36:5:147:U:H5'	118.75	0.48
69:O3:97:SER:OG	36:5:3174:A:OP1	241.26	0.48
1:6:1032:G:H2'	1:6:1033:C:H6	1.77	0.48
1:6:1054:U:H2'	1:6:1055:U:C6	2.45	0.48
1:6:1673:G:H2'	1:6:1674:C:C6	2.48	0.48
1:6:28:A:H2'	1:6:29:U:C6	2.49	0.48
1:6:675:U:H2'	1:6:676:G:H8	1.79	0.48
1:6:882:U:H2'	1:6:883:C:C6	2.49	0.48
13:C1:84:ILE:HG23	13:C1:111:VAL:HG11	1.96	0.48
13:C1:7:VAL:HG13	13:C1:8:GLN:H	1.79	0.48
14:C2:119:SER:OG	14:C2:120:VAL:N	2.47	0.48
15:C3:138:ASN:OD1	15:C3:139:TRP:N	3.34	0.48
18:C6:82:ARG:HH12	18:C6:114:ARG:HB2	3.02	0.48
21:C9:42:GLY:HA2	21:C9:84:LYS:HG3	3.09	0.48
26:D4:10:ARG:NH1	1:6:778:G:N7	431.60	0.48
30:D8:12:VAL:HG22	30:D8:30:VAL:HG12	2.89	0.48
33:E1:95:HIS:CG	33:E1:96:LYS:H	4.04	0.48
39:L2:5:ILE:HG12	39:L2:8:GLN:OE1	2.14	0.48
36:1:2514:U:H5'	45:L8:68:ARG:HG3	1.96	0.48
47:M0:16:PRO:O	47:M0:18:PRO:HD3	2.14	0.48
50:M4:20:VAL:HG22	50:M4:66:THR:OG1	2.13	0.48
55:M9:17:VAL:HG22	55:M9:18:GLY:H	4.55	0.48
56:N0:28:ARG:HE	56:N0:99:ARG:HH21	2.04	0.48
1:2:336:G:H21	10:S8:7:SER:HB2	1.79	0.48
11:S9:159:ALA:HB3	11:S9:162:SER:OG	2.46	0.48
36:1:121:A:C6	45:L8:129:PRO:HG3	2.48	0.48
36:1:1638:A:N3	36:1:1709:C:H1'	2.29	0.48
36:1:2269:U:O2'	36:1:2271:A:N7	2.32	0.48
1:2:1365:C:H5''	18:C6:28:LEU:HD22	1.96	0.48
1:2:1683:C:O2'	1:2:1684:U:O5'	2.31	0.48
37:3:62:U:O4	37:3:63:A:N6	2.47	0.48
38:4:135:G:OP2	61:N5:56:ARG:NH2	2.39	0.48
38:4:21:C:OP1	41:L4:193:LYS:HD2	2.13	0.48
36:5:1594:A:O2'	36:5:1614:C:O2	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
79:Q3:18:TYR:HA	36:5:2131:A:N6	226.75	0.48
36:5:2636:A:H5''	36:5:2637:A:H5'	1.96	0.48
36:5:2716:U:H2'	36:5:2717:U:H5'	1.96	0.48
36:5:3218:A:C8	36:5:3277:U:H5'	2.48	0.48
36:5:359:U:H2'	36:5:360:G:O4'	2.13	0.48
1:6:1046:G:H1	1:6:1072:C:H42	1.60	0.48
1:6:1216:C:O2'	1:6:1444:A:N1	2.44	0.48
1:6:1696:G:O2'	1:6:1698:G:N7	2.47	0.48
1:6:407:A:H2'	1:6:408:C:C6	2.48	0.48
38:8:102:U:H2'	38:8:103:G:C8	2.49	0.48
38:8:40:A:H2'	38:8:41:A:C8	2.49	0.48
12:C0:49:LEU:HB3	12:C0:55:VAL:HG13	1.95	0.48
16:C4:124:ASP:O	16:C4:125:SER:HB2	2.12	0.48
1:2:901:G:N2	16:C4:54:GLU:OE1	2.47	0.48
18:C6:143:ARG:NH1	1:6:1191:U:H5'	348.66	0.48
19:C7:20:TYR:O	19:C7:24:LEU:HD13	2.14	0.48
21:C9:4:VAL:HG11	21:C9:137:ALA:HA	2.94	0.48
22:D0:39:SER:O	22:D0:43:LYS:HB2	2.69	0.48
26:D4:132:ARG:HG2	26:D4:133:ASN:OD1	6.71	0.48
28:D6:70:LYS:HG2	28:D6:72:HIS:CE1	5.36	0.48
28:D6:82:ARG:O	28:D6:84:VAL:HG12	2.13	0.48
28:D6:84:VAL:HG13	28:D6:85:ARG:N	2.29	0.48
30:D8:52:ASP:N	30:D8:52:ASP:OD2	4.05	0.48
33:E1:121:CYS:H	33:E1:130:VAL:HG11	6.93	0.48
40:L3:370:PHE:CD1	40:L3:376:LYS:HA	2.49	0.48
40:L3:50:LYS:HD3	40:L3:330:GLY:O	2.13	0.48
40:L3:55:THR:O	40:L3:56:ILE:HD12	2.13	0.48
42:L5:105:ILE:O	42:L5:109:THR:HG23	2.13	0.48
44:L7:39:GLU:O	44:L7:43:ILE:HG13	2.14	0.48
44:L7:70:LYS:NZ	36:5:519:A:O5'	315.88	0.48
47:M0:38:LYS:HD3	47:M0:41:ALA:HB2	2.66	0.48
46:L9:47:LYS:NZ	50:M4:5:SER:HB2	2.29	0.48
50:M4:25:LYS:HD2	50:M4:62:GLN:HG2	1.95	0.48
51:M5:60:VAL:HG21	38:8:142:C:H4'	104.56	0.48
50:M4:55:ARG:HD3	56:N0:70:THR:HB	2.38	0.48
36:1:1609:C:H5''	61:N5:125:ARG:HH11	1.79	0.48
68:O2:26:HIS:O	68:O2:28:VAL:N	2.89	0.48
51:M5:143:ARG:NH2	71:O5:92:LEU:HD23	2.21	0.48
4:S2:169:LEU:HD22	4:S2:218:ILE:HG23	1.95	0.48
4:S2:230:TRP:CE2	24:D2:68:ARG:HD2	3.74	0.48
5:S3:143:ARG:HB3	35:SM:109:GLY:HA3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:173:TYR:CG	9:S7:181:ILE:HD11	4.91	0.48
9:S7:30:SER:HB2	9:S7:34:LEU:HB2	2.53	0.48
9:S7:28:GLU:OE1	9:S7:35:LYS:HG2	3.93	0.48
36:1:1054:A:H5''	36:1:2637:A:H61	1.78	0.47
36:1:1654:A:C2'	36:1:1655:G:H5'	2.43	0.47
36:1:2254:U:H2'	36:1:2261:G:N2	2.29	0.47
36:1:230:U:H2'	36:1:231:G:O4'	2.14	0.47
36:1:2393:G:H4'	40:L3:252:ILE:HD13	1.95	0.47
36:1:2941:A:O5'	36:1:2943:G:H4'	2.14	0.47
36:1:3006:A:C2	36:1:3141:A:C4	3.02	0.47
36:1:3210:A:H2'	36:1:3211:C:C6	2.48	0.47
1:2:1487:A:H2'	1:2:1488:G:H8	1.79	0.47
1:2:1488:G:H3'	1:2:1515:A:H61	1.79	0.47
1:2:569:C:H2'	1:2:570:A:O4'	2.14	0.47
42:L5:140:ARG:NH2	36:5:1080:A:OP2	230.30	0.47
36:5:1223:A:OP2	36:5:1285:G:N2	2.43	0.47
36:5:1641:U:O2'	36:5:1642:A:H3'	2.14	0.47
36:5:1856:C:H2'	36:5:1857:C:C6	2.48	0.47
36:5:2890:A:N1	36:5:2913:C:N3	2.61	0.47
36:5:3386:G:H2'	36:5:3387:U:H6	1.79	0.47
36:5:730:C:H2'	36:5:731:U:H6	1.78	0.47
1:6:1509:C:H2'	1:6:1510:U:O4'	2.14	0.47
1:6:15:U:H2'	1:6:16:G:O4'	2.13	0.47
1:6:333:A:C6	1:6:334:G:C6	3.02	0.47
1:6:396:G:N2	1:6:399:A:OP2	2.47	0.47
12:C0:29:GLN:OE1	12:C0:39:ASN:ND2	3.03	0.47
18:C6:55:VAL:HG22	18:C6:59:LYS:HE3	1.95	0.47
25:D3:86:PHE:CZ	25:D3:88:PRO:HA	3.37	0.47
31:D9:9:SER:HA	1:6:1451:C:OP1	411.01	0.47
40:L3:315:GLY:HA2	36:5:3379:C:H4'	214.51	0.47
41:L4:62:ALA:HB3	41:L4:90:PHE:CE2	2.49	0.47
47:M0:80:SER:O	47:M0:84:ALA:HB2	2.14	0.47
47:M0:89:VAL:HG13	47:M0:136:PHE:CE1	2.49	0.47
49:M3:114:GLN:HA	49:M3:117:LYS:HD2	4.15	0.47
53:M7:126:ARG:HH11	53:M7:138:LYS:HB2	1.78	0.47
54:M8:143:PRO:HB2	54:M8:146:SER:OG	2.46	0.47
55:M9:130:ASN:HB3	55:M9:131:ALA:H	1.42	0.47
68:O2:41:VAL:HA	68:O2:46:PHE:HD2	1.79	0.47
69:O3:31:LYS:NZ	69:O3:35:VAL:O	2.37	0.47
72:O6:70:ARG:HG2	72:O6:87:VAL:HG21	1.95	0.47
3:S1:81:PHE:HB2	3:S1:82:ARG:H	1.51	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S2:74:PRO:C	4:S2:76:LEU:H	2.31	0.47
5:S3:101:GLN:HA	5:S3:104:SER:HB3	2.25	0.47
6:S4:252:ARG:NH2	6:S4:252:ARG:HB3	4.53	0.47
6:S4:3:ARG:HB3	1:6:93:A:H1'	326.33	0.47
7:S5:77:TYR:HB3	7:S5:84:LYS:HA	1.96	0.47
11:S9:77:ILE:O	11:S9:81:VAL:HG23	2.51	0.47
35:SM:153:UNK:O	35:SM:155:UNK:N	2.47	0.47
34:SR:258:THR:HB	34:SR:275:ARG:HH12	1.79	0.47
36:1:1266:G:H1	36:1:1275:C:H42	1.62	0.47
36:1:1766:G:H8	36:1:1766:G:OP2	1.97	0.47
36:1:208:C:H2'	36:1:209:A:O4'	2.14	0.47
36:1:2675:C:H41	48:M1:22:SER:HB3	1.79	0.47
36:1:3047:U:O2'	36:1:3048:A:H5'	2.13	0.47
36:1:438:A:C2	36:1:439:C:H1'	2.48	0.47
1:2:1146:G:N3	1:2:1635:A:H2	2.12	0.47
1:2:560:U:H2'	1:2:561:G:C8	2.49	0.47
1:2:561:G:N2	1:2:584:C:O2	2.46	0.47
1:2:973:A:H2'	1:2:974:A:C8	2.49	0.47
36:5:2663:G:H2'	36:5:2664:C:O4'	2.14	0.47
36:5:352:A:H61	36:5:365:A:H5"	1.79	0.47
36:5:696:C:HO2'	36:5:697:A:H8	1.58	0.47
1:6:1491:U:H5'	1:6:1492:A:OP1	2.14	0.47
37:7:89:G:N2	37:7:91:G:H3'	2.29	0.47
84:B:76:8AN:H8	86:B:3401:SPS:H81	1.96	0.47
12:C0:11:ILE:HD11	12:C0:42:VAL:HG22	1.94	0.47
13:C1:38:ALA:HB2	13:C1:60:PHE:CD1	3.85	0.47
14:C2:29:LYS:HE2	14:C2:100:TRP:NE1	2.77	0.47
2:S0:198:MET:SD	19:C7:85:VAL:HG11	2.54	0.47
22:D0:100:VAL:O	22:D0:103:ILE:HG22	2.14	0.47
23:D1:16:LYS:HE3	23:D1:16:LYS:HB2	1.68	0.47
26:D4:10:ARG:HB2	26:D4:24:VAL:HB	2.72	0.47
1:2:1797:A:H5'	28:D6:95:ARG:CZ	2.44	0.47
30:D8:64:ARG:HB3	30:D8:65:ARG:H	2.74	0.47
42:L5:232:ASP:OD2	42:L5:232:ASP:N	2.45	0.47
44:L7:158:LYS:HG2	44:L7:203:TRP:CH2	2.48	0.47
45:L8:170:CYS:O	45:L8:174:GLY:N	2.47	0.47
46:L9:71:VAL:O	46:L9:75:VAL:HG23	2.53	0.47
49:M3:99:HIS:CD2	36:5:156:G:C8	80.83	0.47
53:M7:70:THR:HG21	53:M7:81:ALA:HB3	2.82	0.47
54:M8:126:GLN:O	54:M8:130:ARG:HG3	2.13	0.47
56:N0:26:ARG:HH11	57:N1:150:THR:HG21	2.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1113:A:H5''	77:Q1:6:ARG:HH21	1.79	0.47
78:Q2:16:THR:C	78:Q2:18:ARG:H	2.17	0.47
78:Q2:35:LEU:C	78:Q2:37:ALA:H	2.17	0.47
79:Q3:29:LEU:O	79:Q3:33:GLN:HG2	2.47	0.47
6:S4:35:PRO:HD2	6:S4:83:PRO:HG2	2.27	0.47
8:S6:138:ALA:HB1	8:S6:142:ARG:HH12	4.06	0.47
1:2:127:G:N7	8:S6:202:ARG:NH2	2.63	0.47
10:S8:10:LYS:HG2	13:C1:133:LYS:HE2	2.41	0.47
10:S8:96:LEU:HD13	10:S8:179:CYS:SG	2.54	0.47
11:S9:103:ASP:O	11:S9:107:ARG:HG2	2.26	0.47
11:S9:41:GLU:OE2	11:S9:126:ARG:NH2	2.95	0.47
11:S9:7:THR:HG21	1:6:758:U:OP1	382.39	0.47
5:S3:224:ASP:HA	34:SR:190:ALA:HB2	4.44	0.47
36:1:114:A:H2'	36:1:115:A:O4'	2.14	0.47
36:1:2221:G:N2	36:1:2223:A:H3'	2.29	0.47
36:1:381:U:H2'	36:1:382:U:H6	1.78	0.47
36:1:981:U:HO2'	36:1:982:C:P	2.36	0.47
36:1:996:A:H2'	36:1:997:A:O4'	2.14	0.47
1:2:1609:U:OP2	18:C6:14:LYS:NZ	2.47	0.47
1:2:1621:U:H2'	1:2:1622:G:C8	2.49	0.47
1:2:1694:A:O2'	60:N4:98:PRO:O	2.32	0.47
1:2:140:A:H61	1:2:281:G:H5''	1.79	0.47
1:2:38:C:C2'	1:2:39:A:H5'	2.44	0.47
1:2:883:C:H2'	1:2:884:A:H8	1.78	0.47
68:O2:61:LYS:NZ	36:5:1339:C:OP1	194.04	0.47
36:5:1481:A:O2'	36:5:1482:A:O5'	2.32	0.47
75:O9:2:ALA:N	36:5:1493:G:O6	122.64	0.47
36:5:1651:U:H3	36:5:1804:A:H61	1.62	0.47
36:5:193:C:H2'	36:5:194:U:H6	1.79	0.47
36:5:2406:C:H2'	36:5:2407:C:C6	2.49	0.47
36:5:2213:A:N1	36:5:2429:G:H1'	2.30	0.47
36:5:2987:A:H2'	36:5:2988:C:C6	2.49	0.47
1:6:1236:A:H61	1:6:1249:U:H3	1.61	0.47
1:6:1394:G:H2'	1:6:1395:G:H8	1.79	0.47
15:C3:37:ILE:HG12	15:C3:54:LEU:HD11	2.23	0.47
16:C4:67:VAL:O	16:C4:71:CYS:N	2.94	0.47
17:C5:18:ARG:NH1	20:C8:90:ASN:OD1	4.96	0.47
1:2:1390:U:OP1	19:C7:5:ARG:HD2	2.15	0.47
21:C9:132:LEU:HA	21:C9:135:ILE:HD11	3.29	0.47
21:C9:28:LEU:HA	21:C9:111:ILE:HD11	4.18	0.47
23:D1:38:LYS:HG3	23:D1:51:VAL:HG23	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:D2:77:PRO:HD2	24:D2:79:PHE:CZ	4.12	0.47
39:L2:79:ASN:ND2	39:L2:165:VAL:HG22	2.30	0.47
40:L3:316:GLU:HG3	40:L3:318:LYS:HE3	1.95	0.47
41:L4:28:ALA:HB2	41:L4:276:LEU:HD22	2.39	0.47
42:L5:122:VAL:C	42:L5:124:GLU:H	3.70	0.47
46:L9:112:ILE:N	46:L9:126:VAL:O	2.52	0.47
47:M0:196:PHE:CG	47:M0:197:VAL:N	3.14	0.47
52:M6:88:VAL:O	52:M6:90:HIS:N	2.48	0.47
55:M9:41:ILE:O	55:M9:45:VAL:HG23	2.39	0.47
61:N5:82:LEU:HD12	61:N5:126:LEU:HD21	1.96	0.47
63:N7:110:ALA:O	63:N7:114:VAL:HG23	2.28	0.47
66:O0:66:LYS:H	66:O0:66:LYS:HD2	4.17	0.47
70:O4:74:ARG:HG2	70:O4:75:ALA:H	2.53	0.47
71:O5:34:GLN:HB3	71:O5:38:ARG:HH11	4.12	0.47
74:O8:24:THR:HG23	74:O8:44:LYS:HB2	4.21	0.47
4:S2:113:LEU:HB2	4:S2:215:PHE:CD1	2.50	0.47
5:S3:25:PHE:HD2	5:S3:37:VAL:HG11	2.62	0.47
6:S4:115:THR:HG1	6:S4:118:GLU:H	1.56	0.47
7:S5:62:VAL:HG13	7:S5:89:ILE:HG12	2.11	0.47
8:S6:177:ARG:NH2	1:6:143:G:N7	311.80	0.47
10:S8:61:GLU:O	10:S8:62:THR:OG1	4.99	0.47
11:S9:146:PHE:HZ	1:6:765:G:N1	430.62	0.47
11:S9:5:PRO:HG3	1:6:380:U:C2	368.35	0.47
11:S9:93:LEU:HA	11:S9:96:VAL:HG12	3.32	0.47
1:2:576:G:OP2	35:SM:102:THR:HG21	2.14	0.47
36:1:1033:U:H2'	36:1:1034:U:C6	2.49	0.47
36:1:1109:U:H2'	36:1:1110:U:C6	2.48	0.47
36:1:1562:C:HO2'	36:1:1563:C:C5'	2.27	0.47
36:1:1699:A:H2'	36:1:1700:G:C8	2.49	0.47
36:1:2218:G:H2'	36:1:2219:A:C8	2.49	0.47
36:1:2659:G:H4'	36:1:2751:G:O2'	2.13	0.47
36:1:2837:A:H8	36:1:2837:A:OP2	1.97	0.47
36:1:544:C:H1'	36:1:548:G:H22	1.78	0.47
36:1:61:A:H2'	36:1:62:A:O4'	2.14	0.47
36:1:748:U:H2'	36:1:749:C:C6	2.50	0.47
1:2:1592:A:H2'	1:2:1593:A:C8	2.50	0.47
1:2:67:A:O2'	1:2:69:G:OP1	2.21	0.47
1:2:702:G:O6	1:2:737:A:N6	2.47	0.47
36:5:1114:U:C4	36:5:1115:G:N7	2.82	0.47
36:5:1307:G:C2	36:5:1308:A:C2	3.03	0.47
36:5:2367:A:H2'	36:5:2368:A:C8	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:2773:C:H2'	36:5:2774:C:H6	1.78	0.47
36:5:771:A:H2'	36:5:772:U:O4'	2.15	0.47
1:6:1286:U:O4	1:6:1287:A:N6	2.46	0.47
1:6:1584:G:H22	1:6:1611:A:P	2.37	0.47
1:6:1645:G:H1	1:6:1756[A]:A:N6	2.12	0.47
1:6:718:U:H5'	1:6:719:U:H5	1.80	0.47
15:C3:136:PRO:O	15:C3:138:ASN:N	2.60	0.47
2:S0:88:LYS:NZ	19:C7:82:ASP:OD2	2.38	0.47
23:D1:81:ASN:OD1	23:D1:81:ASN:N	2.90	0.47
29:D7:49:HIS:HE1	29:D7:70:LYS:HG2	3.01	0.47
25:D3:90:ASP:OD2	32:E0:12:GLY:HA2	2.31	0.47
39:L2:149:ARG:HH22	39:L2:155:LYS:HE2	1.77	0.47
39:L2:149:ARG:HH12	39:L2:253:GLN:CB	9.05	0.47
40:L3:347:SER:HB2	40:L3:350:ALA:CB	3.79	0.47
49:M3:14:PHE:O	36:5:798:G:O2'	141.35	0.47
50:M4:36:VAL:HG11	50:M4:55:ARG:NH2	2.29	0.47
51:M5:58:GLY:HA3	51:M5:142:ILE:CD1	2.45	0.47
51:M5:17:ASP:OD2	51:M5:17:ASP:N	2.41	0.47
53:M7:71:ALA:O	53:M7:74:LYS:HB2	3.63	0.47
54:M8:25:TYR:HA	54:M8:28:LEU:HD12	2.30	0.47
57:N1:122:GLN:HB2	57:N1:124:VAL:HG22	7.67	0.47
59:N3:27:ASP:C	59:N3:29:SER:H	2.97	0.47
79:Q3:38:ASP:OD1	79:Q3:45:LYS:HB3	2.13	0.47
4:S2:49:LYS:HE3	4:S2:246:GLU:OE1	3.66	0.47
7:S5:92:ARG:NH2	7:S5:169:ASN:OD1	2.50	0.47
34:SR:179:LYS:NZ	34:SR:191:ASP:OD1	2.30	0.47
36:1:1549:U:H2'	36:1:1550:C:C6	2.49	0.47
36:1:168:U:H2'	36:1:169:U:H6	1.78	0.47
36:1:2593:A:H4'	36:1:2594:C:O5'	2.14	0.47
36:1:2611:U:H2'	36:1:2612:U:C6	2.50	0.47
36:1:563:U:H2'	36:1:564:G:C8	2.48	0.47
36:1:979:U:H1'	36:1:980:A:C4	2.49	0.47
1:2:225:A:H2'	1:2:226:A:O4'	2.15	0.47
1:2:338:C:H1'	10:S8:5:ARG:HB3	1.96	0.47
1:2:777:C:N4	26:D4:10:ARG:HB3	2.30	0.47
1:2:973:A:H2'	1:2:974:A:H8	1.79	0.47
37:3:54:U:H4'	37:3:55:A:O5'	2.14	0.47
36:5:1528:G:H1	36:5:1832:C:N4	2.12	0.47
36:5:18:G:N2	38:8:142:C:C2	2.83	0.47
36:5:27:C:H1'	36:5:328:U:H1'	1.95	0.47
36:5:2943:G:H2'	36:5:2944:U:O4'	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1133:A:H2'	1:6:1134:C:O4'	2.14	0.47
18:C6:9:THR:HA	1:6:1340:U:O4	434.62	0.47
1:6:1503:A:H2'	1:6:1504:G:O4'	2.14	0.47
1:6:546:U:H2'	1:6:547:U:C6	2.50	0.47
38:8:37:A:H5''	38:8:39:G:O4'	2.15	0.47
13:C1:105:LYS:HA	13:C1:105:LYS:HD3	2.29	0.47
24:D2:53:ILE:HG12	24:D2:60:LYS:HB2	1.96	0.47
24:D2:82:LYS:C	24:D2:84:GLY:H	2.18	0.47
26:D4:57:VAL:HB	26:D4:60:PHE:HE2	4.25	0.47
26:D4:8:ARG:HB3	1:6:780:A:O2'	436.94	0.47
15:C3:15:ALA:HB2	29:D7:20:LYS:HG3	2.33	0.47
29:D7:37:CYS:HB2	29:D7:40:CYS:HB2	6.37	0.47
43:L6:43:LEU:HD21	43:L6:85:ILE:HG13	2.00	0.47
49:M3:10:LEU:HD23	54:M8:166:LEU:HD11	1.97	0.47
51:M5:7:LEU:HD23	51:M5:46:ASP:HB3	2.81	0.47
36:1:1507:G:N7	53:M7:129:THR:HG23	2.29	0.47
53:M7:18:ARG:HG2	53:M7:147:GLU:HB3	5.11	0.47
56:N0:77:VAL:HG11	56:N0:106:LEU:CD1	2.44	0.47
56:N0:106:LEU:HD23	56:N0:110:MET:HG2	1.96	0.47
59:N3:79:VAL:HB	59:N3:118:VAL:HG13	1.95	0.47
63:N7:38:PHE:O	63:N7:40:HIS:ND1	2.36	0.47
65:N9:24:PRO:HB2	65:N9:26:THR:HG23	8.20	0.47
67:O1:23:VAL:HB	67:O1:28:ARG:HG2	1.97	0.47
68:O2:19:ARG:HE	68:O2:33:ARG:CB	2.88	0.47
79:Q3:59:CYS:C	79:Q3:61:LYS:H	2.17	0.47
3:S1:51:SER:HB3	3:S1:57:ALA:N	4.79	0.47
36:1:1611:G:H2'	36:1:1612:A:H8	1.78	0.47
36:1:1615:C:H2'	36:1:1616:U:H6	1.80	0.47
36:1:2563:G:OP1	45:L8:27:THR:OG1	2.27	0.47
36:1:2878:G:H2'	36:1:2879:C:C6	2.50	0.47
36:1:549:U:H2'	36:1:550:A:C8	2.49	0.47
36:1:88:A:H61	36:1:98:G:H1'	1.80	0.47
36:1:911:C:N4	39:L2:3:ARG:HD3	2.29	0.47
1:2:229:U:H3	1:2:236:A:N6	2.10	0.47
1:2:780:A:H8	26:D4:8:ARG:HB3	1.78	0.47
36:5:1001:G:N3	36:5:1041:U:H5'	2.30	0.47
36:5:129:U:H2'	36:5:130:A:C8	2.50	0.47
36:5:1627:U:H2'	36:5:1814:A:N6	2.29	0.47
36:5:1856:C:H2'	36:5:1857:C:H6	1.80	0.47
36:5:2726:C:C2	36:5:2728:G:C2	3.03	0.47
36:5:3060:C:H1'	36:5:3332:U:H1'	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:L6:82:ARG:HG3	36:5:500:C:H5''	250.02	0.47
1:6:1112:G:O2'	1:6:1133:A:N6	2.39	0.47
8:S6:4:ASN:ND2	1:6:152:U:O2	307.31	0.47
20:C8:122:HIS:ND1	1:6:1558:U:OP1	362.18	0.47
1:6:197:A:H2'	1:6:198:A:C8	2.50	0.47
1:6:291:G:H2'	1:6:292:U:C5	2.49	0.47
13:C1:55:ASP:HB3	13:C1:58:CYS:HB2	1.99	0.47
15:C3:91:LEU:HB3	15:C3:122:ILE:HG12	2.32	0.47
1:2:1525:A:OP1	21:C9:93:HIS:ND1	2.48	0.47
25:D3:74:VAL:HG11	25:D3:104:LEU:HD11	1.96	0.47
40:L3:44:THR:O	40:L3:340:LYS:HG2	2.70	0.47
40:L3:345:ASN:OD1	40:L3:347:SER:HB2	2.14	0.47
41:L4:120:TYR:HD1	41:L4:120:TYR:O	2.16	0.47
41:L4:36:HIS:HD2	41:L4:242:ALA:HB1	3.35	0.47
41:L4:82:THR:O	41:L4:82:THR:OG1	2.33	0.47
42:L5:158:ARG:HH21	37:7:46:A:P	286.09	0.47
43:L6:102:ASN:ND2	43:L6:104:GLU:HB2	2.29	0.47
44:L7:232:ARG:HD2	44:L7:236:ILE:HA	2.33	0.47
50:M4:116:GLU:HG3	52:M6:197:LEU:HD23	5.02	0.47
51:M5:21:PHE:O	51:M5:25:VAL:HG23	2.15	0.47
36:1:841:A:H5'	55:M9:125:LYS:O	2.15	0.47
55:M9:153:LYS:HA	55:M9:156:ASN:HB2	3.41	0.47
59:N3:58:VAL:HG22	59:N3:76:ALA:HB3	1.98	0.47
59:N3:93:LEU:H	59:N3:93:LEU:HD23	1.79	0.47
63:N7:27:LYS:HB3	63:N7:42:LEU:HD13	4.09	0.47
65:N9:21:ILE:O	65:N9:22:LYS:NZ	8.02	0.47
67:O1:20:LEU:O	67:O1:28:ARG:NH2	2.81	0.47
1:2:1066:C:H1'	3:S1:146:GLN:HG2	1.95	0.47
3:S1:140:ILE:HG22	3:S1:213:ARG:HB2	1.96	0.47
4:S2:121:VAL:O	4:S2:125:ILE:HG13	2.15	0.47
4:S2:88:LYS:HD3	4:S2:89:GLN:H	5.38	0.47
5:S3:135:GLU:HG3	5:S3:153:ALA:HB2	4.77	0.47
5:S3:20:GLU:OE2	5:S3:76:ARG:NH2	3.96	0.47
7:S5:151:GLY:HA3	7:S5:156:ARG:H	4.97	0.47
34:SR:73:LEU:HD23	34:SR:80:ALA:HA	1.96	0.47
36:1:1391:C:C2	68:O2:103:LYS:HD3	2.50	0.47
36:1:308:A:H5'	36:1:2223:A:H1'	1.95	0.47
36:1:2270:A:H2'	36:1:2271:A:C8	2.50	0.47
36:1:307:A:H2'	36:1:308:A:C8	2.49	0.47
36:1:2107:A:C2	36:1:3344:A:H8	2.32	0.47
36:1:339:C:OP1	36:1:1380:G:O2'	2.27	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1119:G:H2'	1:2:1120:U:C6	2.50	0.47
1:2:1140:G:H2'	1:2:1141:G:H8	1.79	0.47
1:2:1237:G:H2'	1:2:1238:A:H8	1.80	0.47
1:2:1352:G:H1	1:2:1373:C:H42	1.62	0.47
1:2:1458:G:H5''	1:2:1459:C:OP2	2.14	0.47
1:2:1469:A:H4'	1:2:1541:G:H4'	1.97	0.47
1:2:180:A:H2'	1:2:181:A:O4'	2.14	0.47
44:L7:209:ASN:HD22	36:5:1333:C:H1'	240.94	0.47
36:5:1566:A:H2'	36:5:1567:U:H5'	1.95	0.47
70:O4:83:ASN:ND2	36:5:1709:C:OP1	215.22	0.47
36:5:191:U:H2'	36:5:192:C:C6	2.49	0.47
36:5:1944:U:H2'	36:5:1945:A:C8	2.49	0.47
36:5:717:C:OP1	36:5:751:A:O2'	2.31	0.47
1:6:1532:U:H2'	1:6:1533:C:O4'	2.15	0.47
1:6:260:U:O2'	1:6:261:U:OP1	2.28	0.47
42:L5:260:PHE:CE2	37:7:121:U:H5'	320.39	0.47
12:C0:27:PHE:CD1	12:C0:40:LEU:HD23	2.50	0.47
12:C0:2:LEU:HD22	1:6:1258:U:H4'	433.08	0.47
17:C5:25:LEU:HD23	17:C5:28:MET:HE2	1.95	0.47
17:C5:28:MET:O	17:C5:32:ASP:HB2	2.15	0.47
20:C8:35:ILE:HB	20:C8:38:VAL:CG1	4.49	0.47
21:C9:34:VAL:HG23	21:C9:53:TRP:HE1	1.78	0.47
22:D0:20:ILE:HD12	22:D0:100:VAL:HG21	3.53	0.47
27:D5:62:VAL:O	27:D5:66:VAL:HG23	2.29	0.47
30:D8:13:ILE:HG23	30:D8:30:VAL:HA	1.95	0.47
40:L3:188:ILE:O	40:L3:192:VAL:HG12	2.15	0.47
41:L4:207:VAL:HB	41:L4:227:THR:HG22	2.32	0.47
45:L8:34:PHE:H	45:L8:39:ALA:HB3	4.23	0.47
36:1:2585:G:N7	45:L8:47:SER:OG	2.48	0.47
47:M0:29:SER:HB2	47:M0:125:LEU:HD12	6.58	0.47
47:M0:52:LEU:HB2	47:M0:152:LEU:HD22	1.96	0.47
36:1:2355:G:H5'	53:M7:139:TYR:CE1	2.50	0.47
54:M8:94:PHE:O	54:M8:96:PHE:N	3.73	0.47
44:L7:121:LYS:CB	57:N1:133:ALA:HB3	2.45	0.47
59:N3:104:ASN:ND2	59:N3:108:GLU:HB2	4.52	0.47
72:O6:89:GLU:O	72:O6:93:ILE:HG12	2.13	0.47
75:O9:28:ARG:HD3	75:O9:36:ARG:O	3.43	0.47
6:S4:86:PHE:HE2	6:S4:102:VAL:HG23	2.61	0.47
8:S6:7:TYR:HD2	8:S6:8:PRO:HD2	1.80	0.47
35:SM:64:LYS:C	35:SM:66:ALA:H	2.11	0.47
5:S3:225:TYR:N	34:SR:189:GLU:O	3.22	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1611:G:H2'	36:1:1612:A:C8	2.49	0.47
36:1:1621:A:H2'	36:1:1622:U:C6	2.50	0.47
36:1:2225:U:H2'	36:1:2226:U:C6	2.50	0.47
36:1:2534:G:H22	36:1:2545:C:N4	2.12	0.47
36:1:2616:C:H2'	36:1:2617:U:H5'	1.95	0.47
36:1:25:U:H4'	36:1:26:A:N7	2.30	0.47
36:1:374:A:HO2'	36:1:376:G:H8	1.62	0.47
1:2:1429:G:C1'	22:D0:74:GLU:HG2	2.45	0.47
1:2:1653:C:N4	1:2:1654:G:C6	2.83	0.47
1:2:434:G:OP1	25:D3:78:LYS:HA	2.15	0.47
1:2:728:U:H2'	1:2:728:U:O2	2.13	0.47
36:5:1018:G:H2'	36:5:1019:G:O4'	2.15	0.47
57:N1:105:PHE:CE2	36:5:1062:A:H4'	244.09	0.47
36:5:2144:A:H1'	36:5:2281:A:N6	2.30	0.47
36:5:2631:U:H4'	36:5:2697:A:C2	2.50	0.47
40:L3:102:LEU:O	36:5:3147:G:H4'	240.89	0.47
36:5:3259:U:H5''	36:5:3261:C:H5	1.80	0.47
36:5:91:G:N7	36:5:93:C:C2	2.82	0.47
1:6:1091:A:H4'	1:6:1092:A:O5'	2.14	0.47
36:5:998:A:O2'	37:7:103:A:N3	2.45	0.47
61:N5:48:SER:OG	38:8:136:G:OP1	84.61	0.47
24:D2:107:SER:HA	1:6:804:A:C8	366.79	0.47
25:D3:53:VAL:HG13	25:D3:72:VAL:HB	1.95	0.47
26:D4:124:ARG:HH11	26:D4:124:ARG:HB3	1.80	0.47
29:D7:35:VAL:HG22	29:D7:79:PHE:HB3	1.96	0.47
33:E1:126:CYS:O	33:E1:128:ALA:N	2.47	0.47
41:L4:186:LYS:O	41:L4:200:THR:N	2.99	0.47
42:L5:290:ILE:O	42:L5:294:ALA:N	3.15	0.47
43:L6:82:ARG:NH1	69:O3:105:SER:O	2.62	0.47
47:M0:99:ILE:HG23	47:M0:123:HIS:CG	4.52	0.47
50:M4:39:ILE:HB	50:M4:43:LYS:HB3	1.96	0.47
54:M8:178:ARG:HD3	64:N8:50:PRO:O	6.47	0.47
57:N1:13:TYR:O	57:N1:16:GLN:HB2	2.91	0.47
59:N3:18:PRO:HA	59:N3:51:ALA:HA	2.03	0.47
60:N4:65:GLU:HA	60:N4:68:ALA:HB3	4.52	0.47
73:O7:17:THR:HG22	73:O7:18:LEU:N	2.29	0.47
78:Q2:59:HIS:O	78:Q2:61:LYS:N	2.41	0.47
79:Q3:88:GLU:HA	79:Q3:91:GLU:HG2	1.95	0.47
79:Q3:87:ARG:O	79:Q3:90:VAL:HG22	5.26	0.47
5:S3:191:ASP:HA	5:S3:192:PRO:HD2	1.81	0.47
6:S4:136:VAL:HG21	6:S4:148:ARG:HH12	3.04	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:31:SER:HB2	9:S7:32:PRO:HD3	1.96	0.47
11:S9:32:GLY:HA3	32:E0:40:TYR:CG	2.50	0.47
34:SR:60:SER:HB3	34:SR:62:LYS:HE3	1.97	0.47
36:1:2747:A:H5'	42:L5:175:HIS:HA	1.96	0.47
1:2:1753:A:C6	1:2:1754:A:C6	3.03	0.47
1:2:340:U:H2'	1:2:341:A:H8	1.79	0.47
36:5:1108:U:H2'	36:5:1109:U:C6	2.50	0.47
36:5:1565:G:H2'	36:5:1566:A:O4'	2.15	0.47
36:5:188:U:H1'	36:5:208:C:H1'	1.97	0.47
36:5:2418:G:H4'	36:5:2419:A:OP2	2.12	0.47
36:5:856:G:C6	36:5:857:G:N1	2.83	0.47
36:5:913:A:H2	36:5:2134:G:N3	2.13	0.47
1:6:1315:U:H2'	1:6:1316:G:O4'	2.15	0.47
1:6:1776:A:H2'	1:6:1777:G:H8	1.80	0.47
1:6:1:U:O2'	1:6:370:A:H5'	2.15	0.47
1:6:481:A:H61	1:6:507:U:H3	1.63	0.47
1:6:718:U:H5'	1:6:719:U:C5	2.50	0.47
37:7:3:U:H2'	37:7:4:U:C6	2.50	0.47
16:C4:127:ARG:HD2	28:D6:22:ARG:HH12	1.78	0.47
16:C4:19:ILE:HG23	16:C4:28:VAL:HG22	1.96	0.47
24:D2:67:GLY:O	24:D2:69:LEU:N	3.53	0.47
25:D3:130:VAL:HG21	25:D3:135:LEU:HD21	1.97	0.47
25:D3:38:PHE:HB3	1:6:359:A:C2	325.19	0.47
24:D2:22:LYS:HZ3	29:D7:3:LEU:H	2.24	0.47
44:L7:156:ILE:O	44:L7:159:GLN:HB2	2.15	0.47
44:L7:39:GLU:O	44:L7:42:ALA:HB3	2.15	0.47
41:L4:330:TYR:CZ	44:L7:49:ALA:HA	2.51	0.47
36:1:3112:G:O2'	46:L9:70:THR:HB	2.15	0.47
47:M0:28:ASP:OD1	47:M0:28:ASP:N	2.86	0.47
49:M3:18:TRP:O	49:M3:20:GLU:N	2.47	0.47
51:M5:85:THR:HA	78:Q2:50:PHE:O	3.81	0.47
41:L4:299:ILE:HD12	54:M8:39:ARG:HB2	3.21	0.47
59:N3:93:LEU:HB2	60:N4:20:LEU:HD22	1.97	0.47
38:4:151:C:C5	61:N5:24:LEU:HD11	2.50	0.47
61:N5:71:THR:O	61:N5:75:LYS:HG2	2.15	0.47
64:N8:84:GLU:O	64:N8:87:ARG:HB2	3.06	0.47
64:N8:128:ARG:HB3	72:O6:8:ALA:CB	4.07	0.47
72:O6:95:ALA:O	72:O6:100:HIS:ND1	2.32	0.47
74:O8:5:ILE:HG23	74:O8:54:LEU:HB2	2.33	0.47
78:Q2:63:LYS:HD3	36:5:2795:U:OP2	213.37	0.47
2:S0:117:GLU:O	4:S2:40:LYS:NZ	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:S4:252:ARG:HH21	6:S4:252:ARG:HB3	4.50	0.47
7:S5:51:VAL:HG13	7:S5:131:GLN:HB2	2.43	0.47
7:S5:58:LEU:HD21	7:S5:167:ARG:NE	3.82	0.47
1:2:68:A:OP2	8:S6:171:LYS:NZ	2.48	0.47
10:S8:172:ARG:O	10:S8:176:SER:N	2.61	0.47
10:S8:20:GLN:NE2	10:S8:22:ARG:O	5.09	0.47
34:SR:110:VAL:HA	34:SR:126:SER:HB2	1.96	0.47
36:1:2232:A:H2'	36:1:2233:A:C8	2.50	0.47
36:1:2539:C:H5'	36:1:2541:U:O4	2.15	0.47
36:1:2746:A:H2	42:L5:146:LEU:HB3	1.79	0.47
1:2:1003:A:H1'	1:2:1005:A:N7	2.30	0.47
1:2:1217:A:H5''	12:C0:44:LYS:HG3	1.96	0.47
1:2:1383:G:OP1	22:D0:89:ARG:NH1	2.45	0.47
1:2:292:U:H2'	1:2:293:U:C6	2.50	0.47
1:2:429:G:OP1	1:2:439:U:H5''	2.15	0.47
1:2:109:G:O2'	1:2:796:A:N1	2.42	0.47
36:5:2570:U:O3'	36:5:2572:C:N4	2.48	0.47
36:5:2707:C:H2'	36:5:2708:C:H6	1.80	0.47
36:5:3288:G:H1'	36:5:3289:G:H5'	1.96	0.47
49:M3:171:ARG:HD3	36:5:770:G:OP1	145.19	0.47
36:5:850:U:H2'	36:5:851:C:C6	2.50	0.47
22:D0:89:ARG:NH2	1:6:1383:G:OP1	446.29	0.47
1:6:1474:G:H2'	1:6:1475:A:H8	1.79	0.47
1:6:834:G:H3'	1:6:835:U:C5	2.50	0.47
12:C0:7:ASP:O	12:C0:11:ILE:HG12	2.15	0.47
25:D3:65:ASN:ND2	25:D3:116:ASP:OD1	4.93	0.47
1:2:531:C:O2'	26:D4:64:PHE:HA	2.15	0.47
39:L2:40:TYR:HB3	39:L2:93:LYS:O	2.14	0.47
40:L3:84:VAL:HG23	40:L3:163:HIS:O	3.49	0.47
40:L3:169:THR:HA	40:L3:170:PRO:HD2	2.21	0.47
40:L3:206:ASP:OD1	40:L3:206:ASP:N	2.45	0.47
42:L5:60:ILE:HB	42:L5:80:SER:HB2	1.97	0.47
46:L9:166:ARG:HH21	46:L9:168:ARG:HH12	12.59	0.47
49:M3:123:ILE:HG22	71:O5:118:ILE:HG12	3.00	0.47
51:M5:36:ILE:HD13	51:M5:106:VAL:HG12	2.58	0.47
51:M5:120:TRP:CE3	36:5:269:G:H5'	132.51	0.47
51:M5:45:PRO:O	51:M5:49:ARG:HB2	4.03	0.47
53:M7:70:THR:HG23	53:M7:73:GLY:H	3.60	0.47
61:N5:135:ILE:O	61:N5:139:ILE:HG22	2.14	0.47
61:N5:48:SER:OG	61:N5:49:LYS:N	3.39	0.47
63:N7:46:ILE:HD12	63:N7:69:LYS:O	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
66:O0:30:THR:HA	66:O0:33:SER:HB3	1.96	0.47
66:O0:40:LYS:HB3	66:O0:101:LEU:HD11	1.97	0.47
67:O1:55:LEU:HD23	67:O1:95:PRO:HB3	1.95	0.47
36:1:1481:A:N1	70:O4:2:ALA:HA	2.30	0.47
5:S3:141:LYS:HE3	5:S3:179:GLN:HG3	1.97	0.47
6:S4:49:ARG:HG3	6:S4:50:ASN:N	4.62	0.47
6:S4:42:LEU:N	6:S4:84:ALA:O	2.40	0.47
11:S9:122:VAL:O	11:S9:125:ALA:HB3	2.14	0.47
34:SR:106:HIS:CE1	34:SR:126:SER:HB3	2.49	0.47
36:1:1157:G:H2'	36:1:1158:A:O4'	2.15	0.47
36:1:1420:C:OP2	41:L4:193:LYS:NZ	2.30	0.47
36:1:1658:G:H2'	36:1:1659:U:H6	1.80	0.47
36:1:1751:G:H5'	74:O8:26:LYS:HZ1	1.79	0.47
36:1:194:U:H2'	36:1:195:U:H6	1.79	0.47
36:1:297:G:N2	36:1:297:G:OP2	2.45	0.47
36:1:790:U:H2'	36:1:791:A:O4'	2.16	0.47
36:1:838:G:H2'	36:1:839:C:O4'	2.15	0.47
1:2:1291:G:C2	1:2:1325:A:C2	3.03	0.47
44:L7:92:ILE:HD11	36:5:1159:A:H5'	227.39	0.47
36:5:1571:A:H1'	36:5:1572:U:C6	2.50	0.47
36:5:1786:G:H2'	36:5:1787:A:C8	2.50	0.47
36:5:2171:G:H2'	36:5:2172:A:H8	1.80	0.47
36:5:23:A:H2'	36:5:24:G:H8	1.80	0.47
77:Q1:15:ARG:NH2	1:6:1126:G:OP1	282.56	0.47
33:E1:133:ALA:HB1	1:6:1251:U:H4'	440.20	0.47
1:6:90:C:O2'	1:6:451:A:H5''	2.14	0.47
1:6:737:A:H2'	1:6:738:G:O4'	2.15	0.47
1:6:831:U:O2'	1:6:832:U:H5'	2.14	0.47
14:C2:97:LEU:HA	14:C2:100:TRP:CE3	2.50	0.47
18:C6:45:ARG:O	18:C6:48:VAL:HG12	2.44	0.47
21:C9:57:ARG:HH11	21:C9:57:ARG:HB2	1.93	0.47
24:D2:7:LEU:HD11	24:D2:37:PHE:CD2	3.61	0.47
25:D3:62:LYS:HE3	25:D3:116:ASP:HA	5.14	0.47
26:D4:80:ALA:HA	26:D4:83:LYS:HB2	3.13	0.47
30:D8:54:LEU:HD12	30:D8:55:VAL:H	4.46	0.47
40:L3:105:VAL:HG21	40:L3:148:LEU:HD13	2.19	0.47
41:L4:22:LEU:HD11	41:L4:26:PHE:HB2	2.15	0.47
42:L5:131:LEU:HA	42:L5:131:LEU:HD13	1.81	0.47
44:L7:166:ASN:HA	44:L7:169:ILE:HD12	3.66	0.47
44:L7:184:LEU:O	44:L7:188:ILE:HG12	3.21	0.47
47:M0:47:PRO:HB3	47:M0:171:TRP:CZ2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:M0:207:GLU:HB3	47:M0:211:ARG:HH22	5.23	0.47
48:M1:101:ASN:HB3	48:M1:130:VAL:HA	1.97	0.47
51:M5:21:PHE:HD2	51:M5:22:LEU:HD13	1.80	0.47
54:M8:153:PHE:O	54:M8:161:LYS:HG2	4.77	0.47
55:M9:132:PHE:CD1	55:M9:138:LEU:HG	2.50	0.47
63:N7:4:PHE:CD2	66:O0:63:SER:HB3	2.50	0.47
65:N9:15:LYS:O	65:N9:18:ARG:HB2	2.15	0.47
70:O4:96:GLU:O	70:O4:99:LYS:HB2	3.63	0.47
71:O5:21:LEU:HD22	71:O5:25:LYS:HG3	4.09	0.47
2:S0:180:GLU:HG3	2:S0:184:LEU:HD23	1.97	0.47
2:S0:76:ILE:HD13	2:S0:98:ILE:HB	1.97	0.47
3:S1:28:GLU:OE2	3:S1:94:LYS:NZ	2.32	0.47
4:S2:238:SER:HA	4:S2:239:PRO:HD2	2.36	0.47
6:S4:148:ARG:HH11	8:S6:201:GLN:HG3	1.80	0.47
7:S5:51:VAL:HG11	7:S5:130:ILE:HG23	2.60	0.47
11:S9:133:HIS:NE2	1:6:513:U:H5'	444.89	0.47
36:1:1129:A:OP1	47:M0:13:LYS:NZ	2.33	0.46
36:1:1145:G:OP1	68:O2:44:ARG:NH1	2.47	0.46
36:1:1488:G:O2'	70:O4:10:ARG:O	2.33	0.46
36:1:2993:G:H2'	36:1:3142:A:N6	2.30	0.46
36:1:3069:G:H2'	36:1:3070:A:H8	1.80	0.46
36:1:595:G:H2'	36:1:596:C:C6	2.50	0.46
1:2:1358:G:H2'	1:2:1359:C:C6	2.50	0.46
1:2:1433:G:H2'	1:2:1434:U:H6	1.80	0.46
1:2:1511:U:H2'	1:2:1512:G:H8	1.77	0.46
1:2:1528:U:H2'	1:2:1529:C:C6	2.50	0.46
44:L7:209:ASN:ND2	36:5:1333:C:H1'	240.96	0.46
36:5:1597:C:H5'	36:5:1696:A:H1'	1.96	0.46
36:5:1715:A:H4'	36:5:1716:U:H3'	1.97	0.46
36:5:2105:G:H2'	36:5:2106:A:C8	2.46	0.46
36:5:2318:U:C4	36:5:2319:U:C4	3.03	0.46
78:Q2:20:HIS:HE1	36:5:2741:C:O2	216.06	0.46
36:5:2350:C:H4'	36:5:3308:C:O2'	2.16	0.46
36:5:547:G:C5	36:5:548:G:H1'	2.50	0.46
1:6:393:C:H4'	1:6:1673:G:O2'	2.15	0.46
1:6:820:U:O2'	1:6:821:U:O5'	2.30	0.46
1:6:862:A:C2	1:6:963:A:C4	3.03	0.46
12:C0:6:GLU:OE2	12:C0:6:GLU:N	5.13	0.46
15:C3:46:THR:OG1	15:C3:49:GLN:HG2	4.43	0.46
18:C6:89:LEU:HG	18:C6:105:LEU:HD23	1.97	0.46
17:C5:119:PHE:HE1	20:C8:121:ALA:HB2	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:L2:117:GLU:OE1	39:L2:163:ARG:NH2	3.19	0.46
39:L2:65:ASP:HA	39:L2:66:PRO:HD3	1.81	0.46
40:L3:227:GLU:HG2	40:L3:270:ARG:HD3	1.98	0.46
40:L3:256:HIS:HA	40:L3:257:PRO:C	2.47	0.46
45:L8:64:ILE:O	45:L8:68:ARG:HG2	2.16	0.46
47:M0:193:ASP:OD2	47:M0:194:GLY:N	2.42	0.46
52:M6:27:LEU:HD21	52:M6:102:LEU:HB2	1.96	0.46
36:1:3308:C:O2	53:M7:69:ARG:HD2	2.16	0.46
55:M9:81:ARG:HG3	55:M9:88:ARG:CZ	2.45	0.46
63:N7:111:LYS:O	63:N7:115:LYS:N	2.68	0.46
65:N9:7:HIS:CG	65:N9:8:THR:N	2.82	0.46
67:O1:19:ARG:HD3	67:O1:35:GLU:HG3	1.96	0.46
70:O4:95:ILE:O	70:O4:99:LYS:HB2	2.15	0.46
8:S6:48:TYR:CD2	8:S6:117:GLY:HA3	2.56	0.46
5:S3:222:VAL:HG21	34:SR:229:LYS:HA	1.96	0.46
36:1:1029:G:H2'	36:1:1030:A:C8	2.50	0.46
36:1:1564:U:H2'	36:1:1565:G:C8	2.50	0.46
1:2:1276:U:OP1	5:S3:146:ARG:HD3	2.15	0.46
1:2:1178:G:C2	1:2:1462:G:C5	3.03	0.46
1:2:18:C:H2'	1:2:19:A:H8	1.80	0.46
1:2:832:U:H2'	1:2:833:U:O4'	2.14	0.46
36:5:1222:G:H1'	36:5:1285:G:N2	2.30	0.46
36:5:270:U:H2'	36:5:271:C:H6	1.80	0.46
36:5:561:C:H2'	36:5:562:C:H6	1.79	0.46
39:L2:181:LYS:HB2	36:5:860:G:C6	212.85	0.46
36:5:929:A:H2'	36:5:930:U:C6	2.51	0.46
1:6:1063:U:H2'	1:6:1064:G:C8	2.50	0.46
12:C0:44:LYS:NZ	1:6:1218:G:OP2	423.82	0.46
1:6:1688:U:H2'	1:6:1689:A:C8	2.50	0.46
1:6:84:A:H2'	1:6:85:A:O4'	2.15	0.46
37:7:62:U:O4	37:7:63:A:N6	2.48	0.46
14:C2:81:ASP:HA	14:C2:82:PRO:HD3	1.80	0.46
17:C5:48:GLY:O	17:C5:50:THR:N	3.05	0.46
24:D2:5:SER:C	24:D2:7:LEU:H	3.17	0.46
30:D8:12:VAL:HB	30:D8:52:ASP:H	1.80	0.46
36:1:2180:G:P	39:L2:174:ARG:HH22	2.37	0.46
40:L3:139:GLN:O	40:L3:141:GLY:N	2.97	0.46
41:L4:292:SER:HG	41:L4:293:SER:H	1.56	0.46
42:L5:34:LYS:O	42:L5:38:THR:HG23	2.15	0.46
37:3:115:G:N2	42:L5:72:ASP:O	2.39	0.46
47:M0:87:LEU:HA	47:M0:138:VAL:HG22	3.61	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:59:PHE:CD2	51:M5:142:ILE:HD11	3.90	0.46
51:M5:164:LEU:HD21	36:5:62:A:H5''	97.92	0.46
36:1:3067:C:H5''	55:M9:58:HIS:CD2	2.50	0.46
64:N8:132:LYS:O	64:N8:136:GLU:HG3	2.39	0.46
71:O5:9:LEU:HB3	71:O5:17:LEU:HD21	2.11	0.46
71:O5:40:SER:OG	71:O5:41:LEU:N	3.12	0.46
74:O8:19:ASP:HB2	74:O8:48:SER:HB2	2.81	0.46
2:S0:101:ARG:HH12	2:S0:104:PRO:HD3	1.80	0.46
2:S0:41:ARG:HG2	2:S0:47:VAL:HG23	3.14	0.46
2:S0:109:ASN:H	4:S2:64:LYS:NZ	3.66	0.46
6:S4:251:GLU:O	6:S4:255:ARG:HG2	4.15	0.46
9:S7:126:LEU:HB2	9:S7:173:TYR:HE2	5.08	0.46
1:2:400:A:H5''	10:S8:25:ARG:HA	1.97	0.46
36:1:1562:C:H2'	36:1:1563:C:C6	2.50	0.46
36:1:2406:C:H2'	36:1:2407:C:C6	2.51	0.46
36:1:3049:A:C2	40:L3:75:ALA:HB2	2.51	0.46
36:1:3278:C:H2'	36:1:3278:C:O2	2.15	0.46
1:2:1503:A:H2'	1:2:1504:G:O4'	2.14	0.46
1:2:1751:C:H2'	1:2:1752:U:O4'	2.16	0.46
1:2:397:A:O3'	10:S8:50:GLY:HA2	2.15	0.46
36:5:1494:U:H4'	36:5:1495:U:O5'	2.16	0.46
36:5:1554:U:H1'	36:5:1555:U:C6	2.51	0.46
36:5:1566:A:H62	36:5:1571:A:H2	1.63	0.46
36:5:2980:U:H2'	36:5:2981:U:C6	2.51	0.46
36:5:3321:C:H2'	36:5:3322:A:H8	1.79	0.46
36:5:595:G:C8	36:5:609:G:C6	3.03	0.46
41:L4:209:TYR:OH	36:5:689:U:O4	87.11	0.46
38:8:72:A:N3	38:8:88:A:O2'	2.48	0.46
24:D2:26:LEU:HD21	24:D2:60:LYS:HD3	1.98	0.46
41:L4:120:TYR:CE2	41:L4:277:PRO:HB3	2.50	0.46
41:L4:299:ILE:HG22	41:L4:300:ARG:O	2.15	0.46
42:L5:107:ARG:HH12	42:L5:120:LYS:HG2	1.80	0.46
42:L5:177:GLU:HG3	42:L5:177:GLU:H	1.45	0.46
42:L5:209:GLU:HG2	42:L5:233:ALA:HB2	1.97	0.46
44:L7:120:THR:OG1	57:N1:132:PRO:HB2	2.15	0.46
47:M0:206:LEU:O	47:M0:210:ILE:N	3.56	0.46
48:M1:92:ARG:HH12	48:M1:94:ARG:HH11	6.60	0.46
50:M4:45:LEU:HD12	50:M4:56:GLN:O	2.45	0.46
51:M5:153:ASP:OD2	51:M5:154:PRO:HD2	2.25	0.46
52:M6:85:ARG:C	52:M6:87:MET:H	2.19	0.46
55:M9:147:ALA:HB1	55:M9:151:ARG:HH21	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:N0:42:TRP:CZ2	56:N0:58:ILE:HG12	3.35	0.46
64:N8:70:LYS:HE2	64:N8:129:PHE:CD2	3.01	0.46
69:O3:93:THR:O	69:O3:95:GLY:N	3.09	0.46
72:O6:58:ILE:HG22	72:O6:90:MET:HG3	1.98	0.46
78:Q2:3:ASN:O	36:5:2655:U:H2'	238.29	0.46
3:S1:157:GLN:H	3:S1:160:HIS:CE1	4.07	0.46
4:S2:205:ARG:HD2	1:6:6:G:OP2	379.19	0.46
4:S2:59:HIS:CE1	4:S2:236:PRO:HB2	3.27	0.46
6:S4:206:ASP:O	6:S4:222:LEU:HB2	2.39	0.46
1:2:339:C:P	10:S8:10:LYS:HZ3	2.36	0.46
10:S8:159:GLN:HB3	10:S8:165:LEU:HD23	1.98	0.46
36:1:1714:A:O4'	36:1:1731:A:N6	2.49	0.46
36:1:2424:A:H2'	36:1:2425:G:O4'	2.15	0.46
36:1:2616:C:C2'	36:1:2617:U:H5'	2.46	0.46
36:1:3090:U:OP1	40:L3:270:ARG:NH2	2.47	0.46
36:1:586:C:H2'	36:1:587:U:O4'	2.16	0.46
36:1:656:A:C2	36:1:1440:G:C2	3.03	0.46
1:2:1087:A:H2'	1:2:1088:A:H8	1.80	0.46
1:2:1283:U:H2'	1:2:1284:C:C5	2.50	0.46
1:2:380:U:O3'	11:S9:2:PRO:HA	2.15	0.46
1:2:446:A:OP1	6:S4:59:ARG:NE	2.23	0.46
37:3:46:A:OP1	42:L5:158:ARG:N	2.47	0.46
37:3:3:U:H2'	37:3:4:U:C6	2.49	0.46
36:5:173:G:HO2'	36:5:174:C:H6	1.61	0.46
36:5:2115:G:H22	36:5:2120:A:H1'	1.81	0.46
36:5:2289:U:H2'	36:5:2290:C:C6	2.50	0.46
36:5:23:A:H2'	36:5:24:G:C8	2.51	0.46
36:5:3163:A:H2'	36:5:3164:C:H5'	1.97	0.46
36:5:33:G:O2'	36:5:51:A:N6	2.46	0.46
36:5:438:A:H4'	36:5:439:C:OP2	2.16	0.46
36:5:543:C:H42	36:5:548:G:H1	1.64	0.46
1:6:1114:G:O2'	1:6:1130:G:O6	2.31	0.46
1:6:1208:A:H8	1:6:1269:U:H6	1.64	0.46
1:6:1469:A:H4'	1:6:1541:G:H4'	1.98	0.46
1:6:231:U:HO2'	1:6:232:U:H6	1.63	0.46
1:6:668:C:H3'	1:6:669:G:N2	2.29	0.46
1:6:653:C:N4	1:6:677:G:H1	2.10	0.46
1:6:686:C:H2'	1:6:687:G:C8	2.50	0.46
1:6:712:G:H2'	1:6:713:A:C8	2.50	0.46
1:6:746:A:H2'	1:6:747:C:O4'	2.15	0.46
1:6:895:G:H2'	1:6:896:U:C6	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:7:113:C:H2'	37:7:114:U:O4'	2.16	0.46
16:C4:16:VAL:HG22	16:C4:33:LEU:HA	1.97	0.46
19:C7:71:PHE:CZ	19:C7:74:GLN:HB2	5.12	0.46
39:L2:29:LEU:HD22	39:L2:101:VAL:HG21	3.69	0.46
40:L3:106:TRP:CG	40:L3:130:PHE:HE1	2.59	0.46
40:L3:280:HIS:HD2	40:L3:328:ILE:HG13	1.80	0.46
40:L3:299:ASP:O	40:L3:301:THR:N	3.07	0.46
45:L8:205:ALA:HA	45:L8:208:GLU:HB2	2.45	0.46
46:L9:49:ASN:C	46:L9:49:ASN:HD22	2.17	0.46
49:M3:167:PHE:HA	49:M3:170:LEU:HD12	3.06	0.46
55:M9:37:SER:HA	36:5:1602:A:H5'	99.42	0.46
36:1:3067:C:OP2	55:M9:62:ARG:NH1	2.48	0.46
56:N0:45:LEU:HD22	56:N0:45:LEU:HA	1.70	0.46
62:N6:37:LYS:H	62:N6:37:LYS:CD	2.23	0.46
63:N7:4:PHE:HD2	66:O0:63:SER:HB3	1.79	0.46
69:O3:29:LEU:HD22	69:O3:75:HIS:CD2	2.51	0.46
71:O5:54:VAL:O	71:O5:58:ILE:HG13	2.18	0.46
36:1:2138:A:C4	73:O7:3:LYS:HB3	2.51	0.46
74:O8:11:PHE:CG	74:O8:54:LEU:HD22	2.51	0.46
2:S0:79:ARG:HH12	2:S0:164:ASN:HB2	1.79	0.46
4:S2:225:LEU:HD13	24:D2:68:ARG:HA	1.98	0.46
5:S3:90:ARG:HH12	5:S3:94:ARG:NH1	11.93	0.46
1:2:788:A:C5	6:S4:19:LEU:HD13	2.50	0.46
6:S4:88:ASP:O	6:S4:100:ARG:HA	2.49	0.46
9:S7:21:ALA:O	9:S7:25:VAL:HG23	2.32	0.46
36:1:1472:U:H2'	36:1:1473:G:C8	2.51	0.46
36:1:2278:C:C2'	36:1:2279:A:H5''	2.45	0.46
36:1:2389:C:H42	36:1:2990:G:H1	1.63	0.46
36:1:279:U:H2'	36:1:280:U:C6	2.50	0.46
36:1:3107:U:O4	36:1:3128:G:N2	2.48	0.46
36:1:653:A:C2	36:1:1443:G:C4	3.04	0.46
1:2:1516:A:OP1	22:D0:88:LYS:NZ	2.32	0.46
1:2:591:A:H2'	1:2:592:A:C8	2.50	0.46
1:2:888:U:O2	1:2:988:A:O2'	2.31	0.46
36:5:106:A:H2'	36:5:107:A:O4'	2.14	0.46
47:M0:119:TRP:CZ3	36:5:1126:G:H5''	256.49	0.46
36:5:1648:A:N6	36:5:1807:G:H1'	2.31	0.46
36:5:2888:U:C6	36:5:2911:A:N6	2.83	0.46
36:5:3046:A:H2'	36:5:3047:U:O4'	2.15	0.46
36:5:3285:C:H3'	36:5:3286:G:H5''	1.98	0.46
1:6:1080:U:O2'	1:6:1081:A:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S3:203:PRO:HB3	1:6:1332:C:H4'	427.43	0.46
1:6:717:C:O2'	1:6:721:U:O4	2.25	0.46
37:7:26:C:H2'	37:7:27:A:O4'	2.15	0.46
20:C8:22:VAL:HA	20:C8:34:THR:HG21	5.59	0.46
1:2:1534:G:C8	27:D5:73:GLY:HA3	2.50	0.46
29:D7:74:SER:O	29:D7:76:GLY:N	2.67	0.46
39:L2:215:ASN:HB2	36:5:2968:G:N7	216.46	0.46
36:1:3140:G:OP1	40:L3:20:LYS:NZ	2.48	0.46
40:L3:305:ILE:HG12	40:L3:321:PHE:CZ	2.50	0.46
41:L4:292:SER:O	41:L4:293:SER:OG	2.31	0.46
48:M1:92:ARG:HH12	48:M1:94:ARG:NH1	6.08	0.46
52:M6:108:ILE:HG13	52:M6:108:ILE:O	4.38	0.46
52:M6:110:PRO:HB2	52:M6:111:PRO:CD	3.09	0.46
53:M7:34:GLN:OE1	36:5:413:U:H5''	154.57	0.46
62:N6:100:HIS:CD2	62:N6:101:PRO:HD2	2.85	0.46
62:N6:55:GLU:HB2	62:N6:108:LYS:HB3	3.53	0.46
54:M8:94:PHE:CZ	64:N8:119:PRO:HD3	2.88	0.46
71:O5:4:VAL:HG21	71:O5:9:LEU:HD11	2.44	0.46
2:S0:189:VAL:HG22	2:S0:190:ASP:H	2.05	0.46
6:S4:22:LYS:HG2	11:S9:4:ALA:HB1	1.97	0.46
9:S7:111:LYS:HB3	9:S7:112:ARG:H	1.92	0.46
9:S7:35:LYS:HZ2	9:S7:36:ALA:H	1.63	0.46
10:S8:8:ARG:NH2	10:S8:28:GLU:OE1	9.21	0.46
6:S4:10:LYS:NZ	11:S9:2:PRO:HB3	3.19	0.46
34:SR:286:GLU:HA	34:SR:287:PRO:HD3	1.82	0.46
34:SR:61:PHE:HB3	34:SR:92:TRP:CD2	2.51	0.46
34:SR:41:THR:HG22	34:SR:62:LYS:HB3	1.96	0.46
36:1:1124:U:O3'	47:M0:15:LYS:NZ	2.47	0.46
36:1:1507:G:N3	36:1:1507:G:H5'	2.30	0.46
36:1:192:C:H2'	36:1:193:C:C6	2.51	0.46
36:1:2343:C:H2'	36:1:2344:U:H6	1.81	0.46
36:1:677:A:H4'	36:1:678:G:O5'	2.15	0.46
36:1:715:A:H5''	64:N8:114:GLY:O	2.15	0.46
36:1:829:U:H3	36:1:895:A:N6	2.14	0.46
1:2:1688:U:H3	1:2:1713:G:H22	1.64	0.46
1:2:328:A:H2'	1:2:329:G:O4'	2.16	0.46
1:2:60:U:O4'	1:2:453:U:H5''	2.16	0.46
46:L9:62:ARG:HD2	36:5:1211:U:P	319.19	0.46
36:5:1189:C:H42	36:5:1315:U:H1'	1.79	0.46
36:5:1313:G:N3	36:5:1318:A:H2	2.14	0.46
36:5:311:C:H42	36:5:2778:G:H1	1.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1218:G:O4'	1:6:1444:A:N6	2.49	0.46
1:6:1238:A:H2'	1:6:1239:U:H5'	1.97	0.46
1:6:1293:U:H3	1:6:1322:A:H61	1.62	0.46
1:6:219:A:N6	1:6:843:U:C2	2.84	0.46
1:6:491:C:N4	1:6:496:G:O6	2.48	0.46
16:C4:52:ARG:HB3	1:6:906:A:OP2	294.09	0.46
1:6:907:A:N1	1:6:1008:G:H1'	2.31	0.46
38:8:71:A:H4'	38:8:72:A:O5'	2.15	0.46
86:B:3401:SPS:H91	86:B:3401:SPS:O1	2.15	0.46
17:C5:67:ALA:O	17:C5:69:GLU:N	2.40	0.46
21:C9:66:TYR:HD2	21:C9:124:ILE:HG12	1.80	0.46
25:D3:55:GLU:HA	25:D3:98:GLU:HG2	1.96	0.46
25:D3:95:PHE:CE1	25:D3:135:LEU:HB3	2.50	0.46
28:D6:15:ARG:HB3	28:D6:16:GLY:H	3.26	0.46
29:D7:22:LYS:HZ3	1:6:864:U:H6	357.32	0.46
41:L4:178:LEU:O	41:L4:182:LEU:HD13	5.72	0.46
42:L5:270:LYS:HG3	42:L5:273:ARG:N	4.95	0.46
42:L5:296:GLN:OE1	42:L5:296:GLN:N	4.94	0.46
43:L6:145:LEU:O	43:L6:149:ILE:N	2.46	0.46
47:M0:147:VAL:O	47:M0:150:GLU:HB3	3.85	0.46
47:M0:56:GLU:HG3	47:M0:161:GLY:HA3	4.71	0.46
47:M0:52:LEU:HD23	47:M0:164:LYS:O	3.68	0.46
36:1:1048:A:H2'	47:M0:22:TYR:CZ	2.51	0.46
47:M0:77:THR:C	47:M0:79:VAL:H	2.18	0.46
51:M5:140:LYS:HB3	51:M5:144:ARG:CZ	3.72	0.46
54:M8:124:LEU:HD23	54:M8:124:LEU:HA	2.18	0.46
54:M8:147:ARG:O	54:M8:150:VAL:HG22	2.16	0.46
67:O1:25:PHE:HD2	67:O1:28:ARG:HD2	1.81	0.46
69:O3:13:HIS:O	69:O3:95:GLY:N	2.99	0.46
70:O4:32:ALA:O	70:O4:33:GLN:HG3	4.16	0.46
74:O8:3:ARG:HH22	36:5:1824:U:H5''	147.59	0.46
79:Q3:10:ILE:HD13	79:Q3:30:GLU:HG2	1.98	0.46
79:Q3:84:ARG:HA	79:Q3:87:ARG:NH2	2.30	0.46
6:S4:148:ARG:HG2	6:S4:148:ARG:H	1.72	0.46
11:S9:176:ASN:OD1	11:S9:179:ARG:NH1	2.49	0.46
36:1:2108:C:H1'	36:1:3344:A:C8	2.50	0.46
36:1:2321:A:H2'	36:1:2322:C:O4'	2.16	0.46
36:1:2883:U:H2'	36:1:2884:C:H6	1.81	0.46
36:1:2943:G:H2'	36:1:2944:U:O4'	2.15	0.46
36:1:3018:C:H2'	36:1:3019:U:O4'	2.16	0.46
36:1:3095:U:H2'	36:1:3096:C:H6	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:12:U:HO2'	1:2:1299:G:HO2'	1.56	0.46
1:2:1318:G:H2'	1:2:1319:A:H8	1.81	0.46
1:2:1332:C:O2'	5:S3:162:GLN:HB3	2.16	0.46
1:2:515:A:N6	1:2:537:G:O2'	2.48	0.46
1:2:947:U:H2'	1:2:948:G:H8	1.81	0.46
36:5:1573:G:C6	36:5:1574:C:H1'	2.50	0.46
36:5:2251:G:C4	36:5:2252:A:C8	3.03	0.46
36:5:3225:C:H2'	36:5:3226:A:C8	2.51	0.46
36:5:501:A:H2'	36:5:502:U:C6	2.51	0.46
36:5:730:C:H2'	36:5:731:U:C6	2.51	0.46
1:6:1687:U:H2'	1:6:1688:U:C6	2.51	0.46
1:6:1689:A:H2'	1:6:1690:G:C8	2.50	0.46
10:S8:142:LYS:NZ	1:6:187:G:N7	276.42	0.46
1:6:649:U:H2'	1:6:650:U:C5	2.51	0.46
1:6:914:G:H8	1:6:914:G:OP2	1.99	0.46
86:B:3401:SPS:H81	86:B:3401:SPS:H71	1.73	0.46
16:C4:103:ARG:O	16:C4:107:ARG:HB2	2.14	0.46
16:C4:80:HIS:ND1	16:C4:113:GLY:O	3.57	0.46
17:C5:18:ARG:HD2	17:C5:36:LEU:O	2.51	0.46
22:D0:37:VAL:HG21	22:D0:109:GLU:HB2	1.96	0.46
23:D1:75:ASN:C	23:D1:77:GLY:H	2.80	0.46
25:D3:79:ASN:H	25:D3:79:ASN:HD22	1.64	0.46
26:D4:62:THR:HA	26:D4:69:SER:HA	2.16	0.46
7:S5:143:ARG:HD3	30:D8:55:VAL:HG11	1.98	0.46
32:E0:18:THR:HG21	1:6:584:C:H1'	389.12	0.46
36:1:3314:A:H5''	40:L3:174:LYS:HD2	1.98	0.46
42:L5:158:ARG:HD3	37:7:46:A:OP1	282.05	0.46
43:L6:42:LEU:HD23	43:L6:84:VAL:HG23	1.98	0.46
44:L7:104:GLN:HA	44:L7:109:THR:HG22	1.95	0.46
44:L7:86:VAL:HG22	44:L7:136:TYR:HB3	3.03	0.46
44:L7:232:ARG:O	44:L7:235:PHE:HB2	2.14	0.46
49:M3:106:GLN:HB3	72:O6:18:THR:OG1	2.81	0.46
51:M5:150:TRP:CH2	51:M5:151:ILE:HG12	2.51	0.46
56:N0:104:GLU:O	56:N0:108:GLN:HG2	2.35	0.46
36:1:1898:G:O2'	59:N3:21:ALA:HB2	2.15	0.46
61:N5:106:ASP:HB3	61:N5:127:THR:OG1	3.99	0.46
61:N5:66:PRO:HG2	71:O5:33:VAL:HG13	2.33	0.46
64:N8:73:LEU:HD13	64:N8:109:TYR:CE1	2.51	0.46
72:O6:42:SER:O	72:O6:46:GLU:HG3	3.59	0.46
76:Q0:94:SER:OG	76:Q0:104:PRO:O	4.06	0.46
6:S4:71:LYS:O	6:S4:90:ILE:HA	2.85	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1039:U:H2'	36:1:1040:A:C8	2.51	0.46
36:1:2516:U:O2	36:1:2594:C:N4	2.48	0.46
36:1:2775:U:H2'	36:1:2776:C:C6	2.51	0.46
36:1:96:G:H5'	49:M3:15:ARG:CZ	2.46	0.46
1:2:1291:G:N2	1:2:1324:G:H22	2.13	0.46
1:2:1459:C:H6	1:2:1459:C:OP2	1.99	0.46
1:2:1490:C:H1'	1:2:1491:U:O4'	2.16	0.46
1:2:1535:U:H5	7:S5:185:ARG:C	2.19	0.46
1:2:538:A:H8	1:2:543:C:H41	1.62	0.46
1:2:589:C:H2'	1:2:590:C:C6	2.50	0.46
1:2:598:U:H2'	1:2:599:A:C8	2.51	0.46
36:5:9:U:C4	36:5:10:C:C4	3.04	0.46
36:5:118:U:C5	36:5:119:U:C4	3.04	0.46
36:5:1627:U:H2'	36:5:1814:A:H62	1.81	0.46
36:5:2267:C:O2'	36:5:2268:U:H5'	2.15	0.46
36:5:2660:G:H5''	36:5:2750:U:O2'	2.16	0.46
54:M8:89:ASP:HB3	36:5:677:A:OP1	133.98	0.46
36:5:956:U:H2'	36:5:957:C:C6	2.51	0.46
1:6:506:A:H3'	1:6:506:A:OP1	2.16	0.46
38:8:97:A:H2'	38:8:98:U:O4'	2.16	0.46
12:C0:52:LYS:HG3	12:C0:54:TYR:CD2	2.51	0.46
13:C1:128:CYS:SG	13:C1:131:ILE:HD11	2.56	0.46
14:C2:67:THR:O	14:C2:69:ALA:N	2.68	0.46
17:C5:53:PRO:HB2	17:C5:57:MET:HE2	3.30	0.46
18:C6:129:PHE:CE1	22:D0:78:THR:HA	2.72	0.46
21:C9:134:ARG:NH1	21:C9:138:GLN:OE1	2.47	0.46
22:D0:64:LYS:HE2	22:D0:66:SER:HB2	7.84	0.46
25:D3:114:LYS:HB3	25:D3:117:ILE:HD11	1.97	0.46
27:D5:91:PRO:HB3	27:D5:101:TYR:HE1	2.59	0.46
1:2:1628:U:OP1	28:D6:88:SER:HA	2.16	0.46
24:D2:57:ARG:NE	29:D7:26:GLN:OE1	2.46	0.46
29:D7:54:VAL:HG12	29:D7:63:LEU:HD12	1.97	0.46
31:D9:36:LEU:HD12	31:D9:38:ILE:HG13	1.98	0.46
40:L3:113:GLU:OE2	40:L3:167:ARG:HB3	4.42	0.46
41:L4:289:ILE:O	41:L4:295:ILE:HG13	2.16	0.46
47:M0:68:ALA:HB1	47:M0:155:ALA:HB1	2.38	0.46
50:M4:102:LYS:HB2	50:M4:102:LYS:HE3	1.57	0.46
52:M6:116:LYS:HG3	52:M6:117:ARG:N	2.31	0.46
53:M7:125:GLN:HB2	53:M7:141:SER:OG	3.24	0.46
56:N0:124:LEU:HD23	57:N1:153:PRO:HB2	2.55	0.46
62:N6:112:ASP:H	62:N6:115:ARG:HB2	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:51:LEU:HB2	63:N7:65:ARG:HD2	2.07	0.46
64:N8:139:ARG:HA	64:N8:143:GLY:O	2.15	0.46
71:O5:34:GLN:O	71:O5:36:LEU:N	2.49	0.46
72:O6:60:LEU:HD13	72:O6:64:SER:HB2	1.98	0.46
74:O8:31:LEU:HA	74:O8:37:PRO:HA	1.96	0.46
75:O9:7:PHE:HB2	36:5:1832:C:O2'	109.06	0.46
79:Q3:19:GLY:HA2	36:5:1925:U:O2	239.51	0.46
2:S0:12:GLU:OE2	2:S0:12:GLU:N	2.46	0.46
2:S0:59:LEU:HD12	23:D1:79:LEU:HD21	7.55	0.46
3:S1:70:LEU:HD21	3:S1:79:HIS:CD2	2.51	0.46
6:S4:42:LEU:HD12	6:S4:109:PHE:CD2	5.62	0.46
8:S6:76:LEU:HD12	1:6:1673:G:H5''	288.42	0.46
8:S6:67:VAL:CG2	8:S6:99:GLY:HA2	2.61	0.46
9:S7:184:GLU:HG2	9:S7:185:ILE:H	3.63	0.46
9:S7:74:GLN:O	9:S7:78:THR:OG1	2.25	0.46
9:S7:94:ALA:O	9:S7:96:ARG:NH1	2.48	0.46
10:S8:106:ALA:HB2	10:S8:165:LEU:HG	2.33	0.46
11:S9:62:ARG:CZ	11:S9:68:LYS:HD3	2.46	0.46
36:1:113:C:H42	36:1:267:G:H1	1.62	0.46
36:1:1385:C:OP2	41:L4:202:ARG:HD3	2.16	0.46
36:1:1462:A:H2'	36:1:1463:U:O4'	2.16	0.46
36:1:1472:U:H5'	55:M9:4:LEU:HB2	1.98	0.46
36:1:196:G:N2	36:1:198:A:H3'	2.31	0.46
36:1:2827:U:O2'	36:1:2829:U:O4	2.20	0.46
36:1:3218:A:H5''	36:1:3219:G:C5	2.51	0.46
36:1:3335:A:N7	36:1:3370:A:O2'	2.43	0.46
36:1:1940:G:N2	36:1:3362:A:H8	2.07	0.46
36:1:383:G:N2	36:1:386:A:OP2	2.42	0.46
36:1:551:A:O2'	36:1:552:G:H8	1.98	0.46
36:1:587:U:C2'	36:1:588:G:H5'	2.46	0.46
1:2:1400:A:H4'	19:C7:60:ARG:NH2	2.30	0.46
1:2:245:U:N3	1:2:248:U:OP2	2.46	0.46
1:2:361:C:H2'	1:2:362:G:C8	2.51	0.46
71:O5:95:PHE:CG	36:5:136:G:H5'	62.25	0.46
36:5:1729:A:H4'	36:5:1730:G:OP2	2.16	0.46
36:5:1807:G:O5'	36:5:1807:G:H8	1.98	0.46
36:5:2566:C:C4	36:5:2567:C:C4	3.04	0.46
36:5:2655:U:H4'	36:5:2656:A:O4'	2.16	0.46
36:5:3063:C:H42	36:5:3080:G:H1	1.64	0.46
36:5:539:C:H42	36:5:552:G:H1	1.62	0.46
1:6:102:U:O4	1:6:361:C:H4'	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S1:149:GLN:HB2	1:6:1066:C:H4'	344.04	0.46
1:6:1572:G:N3	1:6:1572:G:H2'	2.31	0.46
1:6:1748:G:C6	1:6:1749:A:C6	3.04	0.46
13:C1:115:PHE:HD2	13:C1:142:VAL:HG23	1.81	0.46
13:C1:91:LEU:HD23	13:C1:102:LYS:HA	1.98	0.46
20:C8:33:THR:HG21	20:C8:40:ARG:HG3	1.97	0.46
24:D2:27:ILE:HD11	24:D2:61:ILE:HD12	1.97	0.46
26:D4:104:SER:HB3	26:D4:107:GLN:HG3	1.98	0.46
28:D6:38:ARG:HH21	28:D6:83:ILE:HG21	1.80	0.46
23:D1:64:GLU:OE1	29:D7:3:LEU:HB2	2.16	0.46
39:L2:204:MET:HG2	39:L2:208:ASP:HB2	3.79	0.46
41:L4:148:ILE:HA	41:L4:149:PRO:C	2.56	0.46
42:L5:148:ILE:HD11	42:L5:160:PHE:CE1	2.50	0.46
42:L5:78:ALA:HB1	42:L5:104:LEU:HD23	2.03	0.46
44:L7:107:ARG:HD3	36:5:1101:G:H5''	236.55	0.46
44:L7:160:ARG:HG3	44:L7:203:TRP:CG	2.51	0.46
44:L7:173:LEU:HB3	44:L7:178:ILE:HB	2.68	0.46
47:M0:208:ASN:O	47:M0:212:GLU:HB2	3.95	0.46
48:M1:155:THR:OG1	48:M1:158:ASP:HB2	2.64	0.46
49:M3:55:ARG:O	49:M3:115:ARG:NH2	2.48	0.46
52:M6:127:LEU:HD21	56:N0:168:PRO:HB2	2.28	0.46
52:M6:57:PHE:CE2	52:M6:72:HIS:HD2	2.34	0.46
54:M8:69:ARG:HG3	36:5:784:A:N7	158.15	0.46
36:1:1721:U:H3'	55:M9:103:ARG:NH2	2.31	0.46
55:M9:88:ARG:HG2	55:M9:88:ARG:H	3.39	0.46
62:N6:125:LYS:HE2	62:N6:126:LEU:O	2.15	0.46
36:1:1635:G:O6	63:N7:17:ARG:HB2	2.16	0.46
64:N8:47:LYS:NZ	36:5:967:A:OP1	185.58	0.46
63:N7:136:PHE:HE2	70:O4:85:VAL:HG13	1.80	0.46
71:O5:81:ARG:HD3	38:8:38:U:C4	69.51	0.46
74:O8:44:LYS:HB3	74:O8:51:LEU:HD11	4.08	0.46
39:L2:57:PRO:HB3	79:Q3:54:ILE:HG22	5.95	0.46
9:S7:31:SER:HA	9:S7:35:LYS:HB3	3.96	0.46
9:S7:38:LEU:HD23	9:S7:41:LEU:HD12	1.98	0.46
10:S8:13:ALA:HB1	1:6:347:G:O2'	294.51	0.46
11:S9:122:VAL:HG23	11:S9:123:HIS:CD2	3.79	0.46
11:S9:83:VAL:HG23	11:S9:85:VAL:HG23	5.30	0.46
36:1:953:G:O2'	36:1:1116:G:H5'	2.16	0.46
36:1:2333:C:H2'	36:1:2334:U:C6	2.51	0.46
36:1:2862:U:H2'	36:1:2863:G:O4'	2.15	0.46
36:1:713:U:H2'	36:1:714:G:O4'	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1068:C:H2'	1:2:1069:A:C8	2.50	0.46
1:2:1182:U:H2'	1:2:1184:A:OP2	2.15	0.46
1:2:1298:U:O2	4:S2:209:ASN:ND2	2.45	0.46
1:2:130:C:O2'	1:2:131:C:OP1	2.29	0.46
1:2:147:A:H2'	1:2:148:A:O4'	2.16	0.46
1:2:1504:G:H21	1:2:1563:C:H1'	1.81	0.46
1:2:1587:A:OP1	18:C6:136:SER:OG	2.28	0.46
1:2:685:A:HO2'	1:2:686:C:P	2.37	0.46
1:2:804:A:N3	24:D2:105:THR:HG22	2.32	0.46
36:5:2567:C:N3	36:5:2568:C:N4	2.64	0.46
36:5:2704:A:O2'	36:5:2705:A:H5''	2.16	0.46
36:5:2805:G:N3	36:5:2967:A:H2	2.14	0.46
36:5:622:A:H2'	36:5:623:U:O4'	2.15	0.46
1:6:1419:G:H2'	1:6:1420:C:O4'	2.16	0.46
1:6:1531:G:H2'	1:6:1532:U:C6	2.50	0.46
1:6:74:U:H5''	1:6:75:U:OP2	2.16	0.46
26:D4:11:LYS:HD3	1:6:784:C:H42	417.04	0.46
13:C1:46:LYS:O	13:C1:50:GLU:HG2	3.93	0.46
21:C9:57:ARG:O	21:C9:61:VAL:HG23	2.23	0.46
22:D0:28:SER:HB3	22:D0:34:LEU:HD13	5.55	0.46
22:D0:18:GLN:O	22:D0:96:PRO:HA	2.16	0.46
39:L2:97:ASN:HB2	39:L2:100:ASN:ND2	2.69	0.46
40:L3:49:TYR:O	40:L3:80:ASP:N	3.07	0.46
42:L5:107:ARG:O	42:L5:111:GLN:HB2	2.16	0.46
42:L5:208:MET:HG3	42:L5:223:PHE:CE1	2.51	0.46
41:L4:314:LYS:HE2	44:L7:150:LYS:O	2.16	0.46
44:L7:198:ALA:O	44:L7:201:PHE:HB3	2.43	0.46
52:M6:157:GLU:OE1	52:M6:160:ARG:NH1	2.49	0.46
55:M9:10:LEU:HD12	55:M9:10:LEU:HA	2.00	0.46
50:M4:60:LEU:HD13	56:N0:152:LEU:HD11	1.98	0.46
56:N0:71:LYS:O	56:N0:73:LYS:HD3	4.57	0.46
54:M8:182:LYS:HE2	64:N8:55:LYS:O	2.16	0.46
69:O3:45:LEU:HD13	69:O3:45:LEU:HA	4.09	0.46
71:O5:89:ARG:HH11	71:O5:89:ARG:HG2	1.80	0.46
74:O8:14:LEU:HD21	74:O8:52:TYR:CD2	4.68	0.46
74:O8:26:LYS:O	74:O8:41:THR:HA	2.16	0.46
3:S1:101:HIS:HA	3:S1:217:LEU:HD22	2.12	0.46
3:S1:157:GLN:O	3:S1:159:SER:OG	4.48	0.46
4:S2:90:THR:HG23	4:S2:92:ALA:H	2.85	0.46
6:S4:230:GLU:HB2	6:S4:233:LYS:HZ2	4.33	0.46
6:S4:92:LEU:HD13	26:D4:17:LEU:HD11	5.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:S4:94:ALA:HB1	26:D4:16:PRO:HB2	1.98	0.46
7:S5:25:LEU:HB2	7:S5:26:ALA:H	1.60	0.46
8:S6:126:ASP:OD2	8:S6:127:THR:HG22	2.83	0.46
8:S6:127:THR:OG1	8:S6:128:THR:N	2.85	0.46
8:S6:134:GLY:HA3	8:S6:158:ILE:HG13	6.12	0.46
8:S6:67:VAL:HA	1:6:1722:A:H1'	273.97	0.46
10:S8:107:THR:H	10:S8:108:PRO:CD	2.29	0.46
1:2:329:G:H5''	10:S8:98:LYS:HB3	1.97	0.46
35:SM:61:ILE:HD12	35:SM:62:ARG:HG2	1.97	0.46
34:SR:83:ALA:HB1	34:SR:110:VAL:HG12	2.05	0.46
34:SR:5:GLU:HG2	34:SR:317:THR:HG23	5.49	0.46
34:SR:64:HIS:ND1	34:SR:86:ASP:OD1	3.17	0.46
36:1:1317:A:O2'	36:1:1318:A:H3'	2.16	0.45
36:1:1400:G:C2	36:1:1401:A:C8	3.05	0.45
36:1:139:G:H2'	36:1:140:C:O4'	2.16	0.45
36:1:1704:A:HO2'	36:1:1705:U:H5	1.63	0.45
36:1:1861:G:O2'	36:1:3066:U:H5''	2.15	0.45
36:1:3393:U:H2'	36:1:3394:U:H6	1.81	0.45
36:1:703:G:O2'	36:1:787:G:H4'	2.15	0.45
1:2:1396:U:H2'	1:2:1397:U:H6	1.81	0.45
1:2:1606:C:H2'	1:2:1607:G:C8	2.51	0.45
1:2:997:G:H2'	1:2:998:A:O4'	2.17	0.45
38:4:14:C:C4	38:4:15:G:C6	3.04	0.45
36:5:1879:A:H2'	36:5:1879:A:N3	2.32	0.45
36:5:2643:A:H2'	36:5:2645:G:O5'	2.16	0.45
1:6:116:U:H2'	1:6:117:U:C6	2.51	0.45
77:Q1:2:ARG:HD2	1:6:1773:C:OP2	310.76	0.45
1:6:839:U:H2'	1:6:840:U:C6	2.51	0.45
18:C6:58:ASP:C	18:C6:60:PHE:H	2.19	0.45
29:D7:20:LYS:HG2	29:D7:21:LEU:HG	2.33	0.45
30:D8:12:VAL:HA	30:D8:30:VAL:HA	2.97	0.45
40:L3:103:THR:HG22	40:L3:104:THR:H	1.81	0.45
40:L3:27:ALA:HB2	40:L3:220:VAL:HG23	1.98	0.45
41:L4:42:VAL:HG12	41:L4:236:LEU:HD21	2.19	0.45
41:L4:135:VAL:HA	41:L4:245:GLY:O	2.16	0.45
41:L4:328:ASN:H	44:L7:166:ASN:HD21	2.49	0.45
45:L8:157:VAL:O	45:L8:160:ILE:HD13	2.16	0.45
46:L9:106:LYS:O	46:L9:107:ASP:HB2	2.16	0.45
50:M4:92:GLU:CD	50:M4:92:GLU:H	2.17	0.45
52:M6:17:GLY:HA3	36:5:1313:G:O3'	266.12	0.45
54:M8:36:LEU:HD11	54:M8:128:ALA:HB1	2.26	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:M8:32:LEU:O	54:M8:35:PHE:HB3	2.16	0.45
54:M8:36:LEU:O	54:M8:40:THR:OG1	2.17	0.45
55:M9:100:ARG:O	55:M9:104:ARG:HG3	5.08	0.45
56:N0:114:HIS:CE1	36:5:1212:A:H1'	310.98	0.45
60:N4:5:ILE:HD12	60:N4:10:GLY:HA2	1.97	0.45
63:N7:46:ILE:HD11	63:N7:49:TYR:HA	2.14	0.45
36:1:2572:C:O3'	63:N7:61:LYS:NZ	2.50	0.45
78:Q2:12:CYS:SG	78:Q2:17:CYS:HB3	1.89	0.45
78:Q2:4:VAL:HB	78:Q2:93:LEU:HB3	5.28	0.45
4:S2:183:ALA:HB1	4:S2:211:LEU:HD21	1.98	0.45
4:S2:140:ARG:HD3	4:S2:222:TYR:CE1	2.51	0.45
5:S3:204:ASP:HB3	1:6:1331:A:N3	422.98	0.45
6:S4:120:SER:O	6:S4:164:LEU:HB2	2.88	0.45
11:S9:63:ASP:O	11:S9:66:ASP:HB2	2.59	0.45
11:S9:88:GLU:HA	11:S9:91:LYS:HE3	3.87	0.45
35:SM:47:ALA:HB2	36:1:2678:A:C8	2.52	0.45
36:1:1245:A:C3'	36:1:1246:G:H5''	2.43	0.45
36:1:279:U:H2'	36:1:280:U:H6	1.81	0.45
36:1:2984:C:H2'	36:1:2985:C:C6	2.52	0.45
86:1:3401:SPS:H81	86:1:3401:SPS:H71	1.60	0.45
1:2:1207:C:H42	1:2:1456:C:H41	1.63	0.45
1:2:319:U:H1'	1:2:323:A:C4	2.51	0.45
37:3:68:C:H2'	37:3:69:C:H6	1.80	0.45
36:5:1560:G:HO2'	36:5:1561:G:P	2.39	0.45
36:5:2296:A:H2	36:5:2918:G:N3	2.15	0.45
36:5:652:G:OP1	36:5:1436:U:O2'	2.32	0.45
1:6:119:A:H1'	1:6:397:A:C5	2.50	0.45
14:C2:131:ASP:OD1	14:C2:132:GLU:N	2.48	0.45
17:C5:90:ILE:HD11	17:C5:112:LEU:HD21	1.98	0.45
18:C6:82:ARG:NH1	18:C6:114:ARG:HB2	3.69	0.45
20:C8:132:ARG:HB3	20:C8:136:GLN:HG3	1.98	0.45
21:C9:54:PHE:CE2	21:C9:104:VAL:HG22	2.51	0.45
21:C9:57:ARG:HH21	21:C9:80:TYR:HB3	1.81	0.45
22:D0:28:SER:HB2	22:D0:112:VAL:HA	1.98	0.45
24:D2:72:CYS:HB3	24:D2:129:VAL:HG13	1.98	0.45
26:D4:104:SER:OG	26:D4:107:GLN:HB2	4.39	0.45
32:E0:24:THR:N	1:6:587:C:OP1	416.65	0.45
14:C2:54:ARG:HD2	33:E1:129:GLY:HA3	1.98	0.45
39:L2:242:ARG:HH11	39:L2:246:LEU:HD12	5.84	0.45
40:L3:187:SER:O	40:L3:191:LYS:HG3	2.57	0.45
40:L3:322:ILE:HG22	40:L3:324:VAL:HG23	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:L3:35:ASP:OD1	40:L3:36:ASP:N	2.88	0.45
43:L6:105:TYR:CE1	43:L6:134:ARG:HD2	2.51	0.45
45:L8:132:VAL:HG11	45:L8:189:LEU:HD12	1.97	0.45
45:L8:243:GLN:NE2	45:L8:247:ASP:OD1	5.46	0.45
47:M0:69:ARG:HD2	47:M0:70:ILE:HD12	1.97	0.45
48:M1:54:VAL:HB	48:M1:57:PHE:HB2	1.97	0.45
54:M8:184:PHE:CG	36:5:2730:G:H4'	190.64	0.45
55:M9:181:ARG:O	55:M9:185:LEU:HG	2.17	0.45
56:N0:139:TYR:CD2	56:N0:140:VAL:HG23	2.51	0.45
36:1:563:U:OP1	56:N0:71:LYS:NZ	2.49	0.45
59:N3:45:ARG:O	59:N3:48:ARG:HB2	2.68	0.45
63:N7:48:ARG:HB3	63:N7:69:LYS:HB3	1.97	0.45
64:N8:61:PHE:CZ	36:5:283:G:H2'	148.49	0.45
68:O2:122:PRO:O	68:O2:123:LYS:HB2	2.17	0.45
70:O4:54:ILE:HD11	70:O4:78:GLY:HA3	3.45	0.45
71:O5:67:ARG:O	71:O5:71:LYS:HB2	2.17	0.45
73:O7:16:HIS:CD2	36:5:360:G:H4'	129.44	0.45
76:Q0:113:ARG:NH2	36:5:1191:U:OP1	289.21	0.45
2:S0:136:ALA:HB1	2:S0:141:ILE:HB	1.97	0.45
2:S0:22:THR:HG22	2:S0:162:CYS:HB2	6.41	0.45
2:S0:60:ALA:O	2:S0:64:ILE:HG13	2.18	0.45
2:S0:64:ILE:HG21	2:S0:89:PHE:HZ	2.82	0.45
4:S2:121:VAL:HG11	35:SM:117:LEU:HB2	1.98	0.45
4:S2:178:ILE:HA	4:S2:196:VAL:HG12	1.98	0.45
4:S2:226:THR:HG23	4:S2:229:LEU:HD13	1.97	0.45
4:S2:225:LEU:HD22	4:S2:230:TRP:HD1	1.81	0.45
6:S4:100:ARG:HH11	6:S4:236:ILE:HB	1.80	0.45
7:S5:43:PHE:CZ	7:S5:90:ILE:HG21	2.93	0.45
8:S6:25:ARG:HA	8:S6:28:PHE:CD2	3.86	0.45
9:S7:103:SER:O	9:S7:107:ARG:HB3	8.50	0.45
9:S7:45:SER:HB3	9:S7:61:PHE:HD2	1.81	0.45
9:S7:75:THR:OG1	9:S7:76:LYS:N	2.49	0.45
35:SM:22:PRO:HD3	48:M1:35:LYS:HG3	1.98	0.45
34:SR:154:VAL:O	34:SR:155:ARG:HD3	3.26	0.45
34:SR:245:PHE:CE1	34:SR:252:LEU:HD13	2.51	0.45
34:SR:267:PRO:HG2	34:SR:269:TYR:CE1	2.51	0.45
34:SR:9:LEU:HA	34:SR:313:TRP:HA	2.25	0.45
36:1:1307:G:H1'	36:1:1308:A:C8	2.52	0.45
36:1:839:C:H4'	36:1:1724:U:H3'	1.98	0.45
36:1:2218:G:H2'	36:1:2219:A:H8	1.81	0.45
36:1:2392:C:H1'	40:L3:266:ARG:HH21	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:270:U:O2'	36:1:318:A:H1'	2.16	0.45
36:1:3307:A:OP1	40:L3:226:PHE:HB2	2.16	0.45
36:5:1773:C:H2'	36:5:1774:C:C6	2.51	0.45
36:5:2261:G:H21	36:5:2262:A:H61	1.63	0.45
36:5:26:A:H2'	36:5:27:C:C6	2.51	0.45
36:5:677:A:H4'	36:5:678:G:O5'	2.16	0.45
1:6:1102:G:H2'	1:6:1103:U:O4'	2.16	0.45
1:6:1245:G:N1	1:6:1249:U:O4	2.49	0.45
8:S6:13:GLN:NE2	1:6:151:G:H21	309.99	0.45
1:6:1634:C:H4'	1:6:1635:A:OP2	2.17	0.45
1:6:1639:C:H2'	1:6:1640:C:O4'	2.15	0.45
1:6:1713:G:H2'	1:6:1714:A:C8	2.51	0.45
16:C4:51:ASP:HB3	1:6:906:A:OP2	290.19	0.45
1:6:982:U:H2'	1:6:983:A:C8	2.51	0.45
38:8:44:A:H2'	38:8:45:C:H6	1.79	0.45
13:C1:102:LYS:HG3	1:6:632:U:OP1	327.76	0.45
14:C2:68:GLU:C	14:C2:70:ASN:H	2.18	0.45
16:C4:99:GLN:HE22	28:D6:46:GLU:H	7.68	0.45
25:D3:61:SER:HB3	25:D3:69:ARG:HD2	1.98	0.45
45:L8:173:MET:HB3	45:L8:175:VAL:HG23	3.53	0.45
47:M0:61:SER:HG	47:M0:63:GLU:HG2	4.12	0.45
48:M1:21:ILE:HG12	48:M1:125:MET:HB3	4.64	0.45
52:M6:110:PRO:O	52:M6:113:ASP:N	4.86	0.45
53:M7:32:THR:HG21	53:M7:84:PRO:HG2	1.98	0.45
36:1:1364:C:H5''	54:M8:3:ILE:CD1	2.46	0.45
55:M9:109:TYR:HB3	55:M9:115:ILE:HG12	5.06	0.45
55:M9:23:TRP:O	55:M9:50:ILE:HA	2.16	0.45
61:N5:86:VAL:O	61:N5:120:LYS:HB2	2.16	0.45
61:N5:64:GLU:HG2	61:N5:85:GLN:O	5.22	0.45
64:N8:111:LYS:HD3	64:N8:113:LEU:HD21	2.60	0.45
64:N8:74:ASN:CB	64:N8:115:LYS:HB2	2.46	0.45
65:N9:44:LYS:HE3	65:N9:44:LYS:HB2	2.97	0.45
70:O4:16:ARG:HH11	70:O4:37:LYS:HD2	1.82	0.45
4:S2:152:HIS:N	4:S2:152:HIS:CD2	3.60	0.45
7:S5:39:GLU:HG2	7:S5:40:ILE:HG22	5.43	0.45
7:S5:89:ILE:HG13	7:S5:89:ILE:H	1.47	0.45
8:S6:88:ARG:HB3	8:S6:91:GLU:HB2	1.98	0.45
34:SR:116:ASP:HB3	34:SR:121:MET:HB3	1.97	0.45
36:1:1194:G:H2'	36:1:1195:A:C8	2.51	0.45
36:1:1818:U:H2'	36:1:1819:U:O4'	2.16	0.45
36:1:2700:G:O2'	36:1:2705:A:N1	2.33	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:3035:A:C4	36:1:3036:G:C8	3.04	0.45
36:1:67:A:O2'	36:1:315:C:O2	2.26	0.45
36:1:353:G:O2'	36:1:354:U:OP2	2.34	0.45
36:1:824:C:H2'	36:1:825:U:C6	2.51	0.45
1:2:110:U:O2'	1:2:797:G:H1'	2.17	0.45
1:2:1291:G:N2	1:2:1324:G:N2	2.64	0.45
1:2:190:C:O2'	1:2:191:C:H5'	2.15	0.45
1:2:252:U:H4'	6:S4:131:LEU:HD12	1.97	0.45
1:2:560:U:H2'	1:2:561:G:H8	1.81	0.45
1:2:647:G:N2	1:2:687:G:H22	2.13	0.45
1:2:992:A:H2'	1:2:993:A:H5'	1.97	0.45
38:4:64:U:H3	38:4:96:A:H61	1.63	0.45
36:5:109:A:N1	36:5:322:U:O2'	2.44	0.45
36:5:1258:U:H2'	36:5:1260:A:OP2	2.16	0.45
68:O2:101:SER:HB3	36:5:1389:G:H5''	128.39	0.45
36:5:2884:C:H2'	36:5:2885:C:H6	1.82	0.45
46:L9:40:HIS:ND1	36:5:3124:G:H5'	311.12	0.45
36:5:3218:A:H5''	36:5:3219:G:C8	2.51	0.45
36:5:3390:G:C2	36:5:3391:A:C8	3.04	0.45
36:5:630:A:H2'	36:5:631:U:C6	2.52	0.45
1:6:1233:G:H1	1:6:1252:C:H42	1.64	0.45
19:C7:44:LYS:NZ	1:6:1387:G:N7	434.78	0.45
1:6:1642:G:H2'	1:6:1643:U:C6	2.51	0.45
1:6:386:G:C6	1:6:387:A:N6	2.84	0.45
1:6:74:U:N3	1:6:76:A:H5''	2.31	0.45
15:C3:129:TYR:HB2	15:C3:135:LEU:HD12	1.99	0.45
18:C6:122:ARG:HG2	1:6:1584:G:H5''	395.11	0.45
24:D2:36:LYS:O	24:D2:40:VAL:HG23	2.27	0.45
1:2:458:G:H5''	26:D4:109:LYS:HZ1	1.82	0.45
26:D4:124:ARG:HD2	1:6:149:C:OP2	327.69	0.45
39:L2:29:LEU:HA	39:L2:76:PHE:CE1	2.37	0.45
40:L3:239:PRO:O	40:L3:242:THR:HG23	2.17	0.45
40:L3:261:MET:HE3	40:L3:261:MET:HB3	2.76	0.45
40:L3:282:ILE:HG21	40:L3:285:VAL:HG22	1.98	0.45
40:L3:37:ARG:CA	40:L3:186:GLY:HA2	3.70	0.45
41:L4:110:ASN:HB3	51:M5:202:TYR:CE2	3.31	0.45
45:L8:111:LYS:HD3	45:L8:111:LYS:HA	1.77	0.45
46:L9:166:ARG:HD2	46:L9:168:ARG:NH1	14.12	0.45
46:L9:47:LYS:HZ2	50:M4:5:SER:HB2	1.81	0.45
48:M1:114:ILE:HG22	48:M1:115:LYS:O	2.36	0.45
50:M4:17:VAL:HG22	50:M4:36:VAL:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:11:GLN:HG2	51:M5:44:ARG:NH2	2.52	0.45
53:M7:32:THR:HG22	53:M7:58:ILE:HG13	1.98	0.45
53:M7:65:SER:O	53:M7:66:SER:HB2	2.16	0.45
55:M9:90:PRO:HG2	55:M9:93:VAL:CG2	3.43	0.45
57:N1:44:ALA:HB2	57:N1:53:PRO:HG2	2.32	0.45
4:S2:243:TYR:HB3	4:S2:246:GLU:HB2	1.97	0.45
5:S3:92:GLN:NE2	5:S3:92:GLN:O	2.49	0.45
8:S6:137:ARG:HB2	8:S6:140:ASN:HB2	3.53	0.45
9:S7:147:ASN:N	9:S7:147:ASN:OD1	2.48	0.45
1:2:333:A:C8	10:S8:49:ARG:HD2	2.51	0.45
11:S9:168:ARG:CZ	11:S9:174:ARG:HH11	7.89	0.45
11:S9:54:ARG:HE	11:S9:54:ARG:HB3	2.84	0.45
34:SR:86:ASP:O	34:SR:88:THR:HG23	2.17	0.45
36:1:1002:A:H2	36:1:1050:U:HO2'	1.60	0.45
36:1:1063:G:H2'	36:1:1097:G:N2	2.32	0.45
36:1:2219:A:H2'	36:1:2220:A:C8	2.52	0.45
36:1:3193:C:H2'	36:1:3194:C:O4'	2.16	0.45
36:1:586:C:OP1	69:O3:70:LYS:NZ	2.38	0.45
36:1:816:A:H5''	36:1:920:A:H62	1.81	0.45
1:2:1471:A:C5	1:2:1472:C:H5	2.35	0.45
48:M1:97:SER:HB3	36:5:2672:G:H1'	328.41	0.45
36:5:298:U:H5''	36:5:299:G:H5'	1.98	0.45
40:L3:329:PRO:HA	36:5:3047:U:H5'	233.64	0.45
40:L3:130:PHE:CZ	36:5:3149:G:H4'	221.23	0.45
1:6:1150:G:N2	1:6:1768:G:H2'	2.32	0.45
1:6:1273:G:N7	1:6:1430:U:H3'	2.32	0.45
1:6:1505:A:C5	1:6:1506:G:H1'	2.52	0.45
1:6:289:U:H2'	1:6:290:G:O4'	2.15	0.45
1:6:393:C:H2'	1:6:394:C:C6	2.52	0.45
1:6:808:U:H2'	1:6:809:A:C8	2.52	0.45
16:C4:42:VAL:HB	16:C4:63:ALA:HB1	4.22	0.45
17:C5:98:ASN:HB3	17:C5:120:SER:OG	2.17	0.45
24:D2:105:THR:HG23	24:D2:110:ILE:HG12	2.05	0.45
26:D4:23:PHE:CZ	26:D4:75:VAL:HG23	6.29	0.45
39:L2:15:ILE:HA	39:L2:15:ILE:HD12	4.76	0.45
39:L2:113:VAL:HG12	39:L2:166:ILE:HA	2.36	0.45
40:L3:117:ARG:CZ	40:L3:175:LYS:HG2	2.83	0.45
40:L3:85:VAL:HG12	40:L3:165:GLN:HE21	2.13	0.45
40:L3:70:ARG:NH1	40:L3:70:ARG:HB3	4.73	0.45
36:1:577:C:H1'	41:L4:340:GLY:O	2.17	0.45
50:M4:20:VAL:HG13	50:M4:68:LEU:O	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:16:SER:OG	51:M5:19:LEU:HB2	2.41	0.45
52:M6:116:LYS:NZ	56:N0:165:TYR:O	2.49	0.45
52:M6:171:LYS:O	52:M6:175:THR:HG22	3.20	0.45
36:1:3095:U:OP1	59:N3:86:ARG:HD2	2.17	0.45
64:N8:133:LEU:HD23	64:N8:133:LEU:HA	1.81	0.45
36:1:2818:U:P	65:N9:2:ALA:HB2	2.57	0.45
2:S0:30:GLN:NE2	2:S0:33:GLN:HG2	10.70	0.45
6:S4:163:ASP:HB3	6:S4:164:LEU:H	3.24	0.45
6:S4:158:ASP:OD2	6:S4:174:LYS:NZ	2.50	0.45
7:S5:73:THR:HG21	18:C6:114:ARG:HE	6.74	0.45
8:S6:70:PRO:O	8:S6:98:ARG:NH1	3.19	0.45
9:S7:15:GLU:O	9:S7:19:GLN:HG3	2.30	0.45
9:S7:166:LEU:HD22	9:S7:166:LEU:H	1.81	0.45
34:SR:242:SER:HB3	34:SR:292:LEU:HD23	1.99	0.45
36:1:1047:A:N3	36:1:2633:U:O2'	2.46	0.45
36:1:2393:G:O2'	36:1:2394:G:OP2	2.27	0.45
36:1:239:G:HO2'	36:1:240:U:P	2.38	0.45
36:1:23:A:H2'	36:1:24:G:O4'	2.17	0.45
36:1:860:G:OP2	39:L2:181:LYS:NZ	2.45	0.45
1:2:1619:C:H1'	30:D8:22:ARG:HE	1.81	0.45
1:2:1629:G:H2'	1:2:1630:U:C6	2.51	0.45
1:2:346:G:H5'	13:C1:79:LYS:HD2	1.97	0.45
37:3:22:A:C6	37:3:23:A:C6	3.05	0.45
36:5:1346:G:H1	36:5:1358:C:H42	1.65	0.45
36:5:1742:U:H2'	36:5:1743:G:C8	2.51	0.45
36:5:193:C:C2	36:5:203:G:C2	3.05	0.45
36:5:2933:A:OP1	36:5:3015:G:H4'	2.17	0.45
36:5:94:G:H2'	36:5:95:A:C8	2.52	0.45
15:C3:102:LEU:HA	15:C3:102:LEU:HD23	1.89	0.45
15:C3:46:THR:HG23	15:C3:49:GLN:OE1	3.21	0.45
15:C3:50:ILE:HG22	15:C3:71:ILE:HD13	1.99	0.45
15:C3:73:ARG:HD3	1:6:859:A:C5	330.93	0.45
1:2:1390:U:H6	19:C7:3:ARG:HG2	1.82	0.45
22:D0:22:ILE:HG22	22:D0:93:LEU:HB2	1.99	0.45
23:D1:8:LEU:HD22	23:D1:8:LEU:HA	2.09	0.45
24:D2:94:LEU:HA	24:D2:95:PRO:HD3	1.86	0.45
30:D8:18:ARG:HD3	1:6:1616:G:H4'	359.94	0.45
39:L2:182:ALA:HB2	36:5:2148:U:O2'	211.04	0.45
39:L2:96:LEU:HG	39:L2:107:VAL:HG12	4.51	0.45
40:L3:223:GLY:HA2	40:L3:271:GLY:HA3	1.98	0.45
36:1:2746:A:OP1	42:L5:179:ARG:NH1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:L5:211:LEU:HD11	42:L5:218:ARG:HB2	1.99	0.45
44:L7:97:PRO:HG3	36:5:1139:G:OP1	230.78	0.45
47:M0:191:LYS:HG2	47:M0:198:LYS:HB2	1.99	0.45
51:M5:93:LYS:HG3	36:5:289:A:C2	146.36	0.45
53:M7:29:THR:OG1	53:M7:29:THR:O	2.30	0.45
55:M9:144:GLN:HA	55:M9:147:ALA:HB3	1.98	0.45
56:N0:99:ARG:NH1	56:N0:126:VAL:O	2.42	0.45
56:N0:36:ILE:O	56:N0:40:ARG:HG3	4.07	0.45
60:N4:23:ARG:NH2	60:N4:25:ASP:OD2	2.46	0.45
68:O2:75:LEU:HA	68:O2:75:LEU:HD23	1.75	0.45
74:O8:45:VAL:O	74:O8:47:GLY:N	3.21	0.45
77:Q1:22:ALA:HA	77:Q1:25:LYS:HB2	1.97	0.45
78:Q2:17:CYS:O	78:Q2:19:LYS:N	2.45	0.45
78:Q2:28:TYR:CB	78:Q2:69:VAL:HB	2.46	0.45
39:L2:178:PRO:HD2	79:Q3:26:VAL:HG23	1.98	0.45
2:S0:179:ARG:HG2	2:S0:183:ARG:HD2	1.99	0.45
3:S1:24:PHE:HA	3:S1:27:LYS:HG2	5.12	0.45
4:S2:38:VAL:O	4:S2:39:THR:OG1	2.27	0.45
6:S4:202:ASP:HB3	13:C1:40:LEU:HD12	3.74	0.45
7:S5:190:ILE:O	7:S5:194:LEU:HB2	2.30	0.45
7:S5:41:LYS:HB2	7:S5:67:PRO:HB2	5.20	0.45
9:S7:31:SER:HA	9:S7:35:LYS:CB	4.21	0.45
10:S8:138:ASN:HB3	10:S8:141:ARG:NH1	4.81	0.45
10:S8:117:TYR:CD1	10:S8:150:ALA:HB2	2.52	0.45
11:S9:155:HIS:O	11:S9:157:ASP:N	2.71	0.45
34:SR:215:GLY:O	34:SR:239:GLU:HG3	2.16	0.45
34:SR:256:THR:N	34:SR:259:GLY:O	2.48	0.45
36:1:1046:A:H2'	36:1:1049:C:C5	2.51	0.45
36:1:1266:G:H1	36:1:1275:C:N4	2.14	0.45
36:1:192:C:H2'	36:1:193:C:H6	1.82	0.45
36:1:2631:U:OP2	57:N1:4:SER:OG	2.35	0.45
36:1:3194:C:N4	36:1:3196:U:O2	2.49	0.45
36:1:650:C:O5'	36:1:650:C:H6	2.00	0.45
36:1:664:U:H2'	36:1:665:A:C8	2.52	0.45
1:2:1530:C:P	27:D5:95:HIS:HB2	2.57	0.45
1:2:190:C:H1'	1:2:191:C:H5'	1.97	0.45
1:2:212:U:C2	1:2:254:A:C2	3.05	0.45
1:2:952:A:H5'	15:C3:98:VAL:HG22	1.98	0.45
1:2:953:G:H2'	1:2:954:G:C8	2.52	0.45
36:5:1807:G:C6	36:5:1808:G:N1	2.85	0.45
36:5:886:C:H4'	36:5:1850:A:C2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:617:U:H5'	1:6:1031:U:O4'	2.17	0.45
1:6:870:C:H2'	1:6:871:G:C8	2.52	0.45
16:C4:122:PRO:HB3	1:6:887:A:H1'	283.38	0.45
1:6:896:U:H2'	1:6:897:C:C6	2.51	0.45
12:C0:24:LYS:HD3	12:C0:63:TYR:CZ	3.38	0.45
16:C4:136:ARG:NH1	1:6:1785:U:OP1	297.95	0.45
20:C8:91:ASP:HB3	20:C8:95:GLY:H	2.06	0.45
26:D4:20:ARG:HB3	26:D4:76:TYR:CD2	2.69	0.45
27:D5:38:HIS:HB3	27:D5:39:ALA:H	3.49	0.45
41:L4:201:GLN:HG3	41:L4:202:ARG:O	2.17	0.45
42:L5:144:VAL:HG12	42:L5:173:VAL:HG22	2.17	0.45
43:L6:158:TYR:CE1	50:M4:115:PHE:HA	2.58	0.45
43:L6:22:ARG:O	43:L6:23:LYS:HD3	2.17	0.45
43:L6:98:VAL:HA	43:L6:101:PHE:CD2	2.52	0.45
36:1:2645:G:P	47:M0:117:GLY:HA2	2.56	0.45
47:M0:38:LYS:HG2	47:M0:41:ALA:HB2	1.99	0.45
47:M0:82:ARG:O	47:M0:82:ARG:HG2	4.46	0.45
48:M1:23:VAL:HG12	48:M1:25:GLU:H	2.96	0.45
52:M6:14:HIS:HE1	52:M6:119:VAL:HG12	1.81	0.45
53:M7:4:TYR:CE1	53:M7:16:SER:HB2	3.29	0.45
60:N4:21:PHE:HB3	60:N4:29:PHE:HB2	1.98	0.45
64:N8:21:ARG:NH1	36:5:1369:A:OP1	183.26	0.45
64:N8:29:PRO:C	64:N8:31:GLY:H	2.64	0.45
70:O4:57:LEU:HB3	70:O4:61:GLN:HB2	1.99	0.45
51:M5:15:GLN:HB3	72:O6:52:PRO:HD2	1.98	0.45
72:O6:53:TYR:HB2	72:O6:76:ARG:HD2	1.97	0.45
3:S1:158:SER:HB2	1:6:875:G:OP1	315.09	0.45
4:S2:102:VAL:HG11	4:S2:129:ILE:HG12	1.98	0.45
7:S5:182:ALA:O	7:S5:186:ASN:ND2	2.49	0.45
11:S9:159:ALA:O	11:S9:165:GLY:HA3	2.80	0.45
34:SR:69:GLN:HG2	34:SR:111:MET:SD	2.57	0.45
36:1:1429:G:C5	41:L4:99:MET:HE1	2.51	0.45
36:1:1506:A:H1'	36:1:1848:G:O6	2.17	0.45
36:1:194:U:H2'	36:1:195:U:C6	2.50	0.45
36:1:2335:G:N2	36:1:2339:C:O2	2.40	0.45
36:1:3326:G:H2'	36:1:3327:G:C8	2.51	0.45
36:1:728:G:H4'	54:M8:47:VAL:HG21	1.99	0.45
1:2:1174:C:H2'	1:2:1175:U:O4'	2.17	0.45
1:2:1353:U:H2'	1:2:1354:G:H5"	1.98	0.45
1:2:226:A:H2'	1:2:227:U:H5'	1.99	0.45
1:2:542:A:H5"	1:2:544:A:C8	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:716:C:N4	1:2:723:G:H21	2.15	0.45
38:4:83:C:H1'	38:4:85:G:N2	2.32	0.45
56:N0:137:ARG:HH12	36:5:1213:G:H5''	324.55	0.45
36:5:138:U:H2'	36:5:139:G:H8	1.78	0.45
36:5:3072:C:H2'	36:5:3073:A:O4'	2.17	0.45
1:6:1209:C:N4	1:6:1454:G:H1	2.15	0.45
1:6:1511:U:H2'	1:6:1512:G:H8	1.81	0.45
1:6:716:C:H1'	1:6:723:G:N2	2.32	0.45
1:6:778:G:C2	1:6:780:A:H5'	2.51	0.45
61:N5:38:LEU:HD12	38:8:147:U:O4'	121.96	0.45
14:C2:52:LEU:HD13	14:C2:85:LYS:NZ	2.49	0.45
7:S5:112:ARG:HH12	18:C6:42:GLU:HG3	4.84	0.45
22:D0:50:LEU:HD22	22:D0:94:GLU:O	4.25	0.45
24:D2:76:SER:HB3	24:D2:77:PRO:HD3	1.98	0.45
26:D4:112:LYS:O	26:D4:116:LYS:HG3	2.17	0.45
26:D4:43:LYS:O	26:D4:47:VAL:HG23	2.17	0.45
26:D4:91:LEU:O	26:D4:96:LEU:N	2.97	0.45
39:L2:43:GLY:O	39:L2:88:ILE:N	2.62	0.45
40:L3:37:ARG:HG2	40:L3:187:SER:N	3.77	0.45
41:L4:143:GLU:HB3	41:L4:144:LYS:NZ	10.32	0.45
41:L4:168:ALA:O	41:L4:172:VAL:HG23	5.08	0.45
41:L4:33:ASP:OD1	41:L4:33:ASP:N	2.62	0.45
42:L5:286:VAL:O	42:L5:290:ILE:HG13	3.52	0.45
44:L7:111:ILE:O	44:L7:112:ASN:HB2	2.17	0.45
44:L7:89:ILE:HD12	44:L7:214:TRP:CH2	2.51	0.45
46:L9:90:MET:HE1	46:L9:179:ILE:O	4.65	0.45
46:L9:9:GLN:O	46:L9:72:LYS:NZ	3.55	0.45
47:M0:68:ALA:HB2	47:M0:158:LYS:HB2	1.99	0.45
49:M3:119:TYR:HD1	49:M3:145:PHE:CE2	2.35	0.45
36:1:950:G:OP1	54:M8:8:LYS:HE2	2.16	0.45
56:N0:135:VAL:HG21	56:N0:144:LEU:HD11	1.99	0.45
61:N5:24:LEU:HB3	61:N5:25:LYS:H	1.43	0.45
70:O4:58:ARG:HG3	70:O4:58:ARG:HH11	1.81	0.45
71:O5:12:LYS:HB2	71:O5:17:LEU:HG	1.98	0.45
77:Q1:1:MET:HE2	77:Q1:5:TRP:HB2	2.33	0.45
77:Q1:22:ALA:O	77:Q1:25:LYS:N	3.38	0.45
3:S1:116:LYS:HG2	3:S1:117:TRP:H	4.45	0.45
3:S1:180:THR:HG23	3:S1:183:GLN:NE2	9.66	0.45
7:S5:30:PRO:HB2	7:S5:33:VAL:HB	1.98	0.45
9:S7:90:VAL:O	9:S7:165:LYS:NZ	4.98	0.45
34:SR:144:LEU:HD21	34:SR:186:PHE:HB3	3.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1567:U:H2'	36:1:1568:U:H4'	1.99	0.45
36:1:3230:G:H4'	50:M4:132:LYS:HD3	1.98	0.45
1:2:1157:A:H2'	1:2:1160:A:N7	2.31	0.45
1:2:1561:U:H2'	1:2:1562:G:C8	2.44	0.45
1:2:448:C:OP1	6:S4:28:ALA:HA	2.16	0.45
1:2:765:G:O6	11:S9:82:ARG:NH1	2.49	0.45
1:2:823:G:H2'	1:2:824:G:C8	2.52	0.45
36:5:151:A:H2'	36:5:152:U:O4'	2.17	0.45
55:M9:110:ARG:NH1	36:5:1720:U:OP2	233.05	0.45
36:5:1495:U:H2'	36:5:1842:A:C2	2.51	0.45
36:5:2960:C:H2'	36:5:2961:G:C8	2.52	0.45
36:5:3225:C:H2'	36:5:3226:A:H8	1.82	0.45
36:5:540:U:H3	36:5:551:A:H61	1.64	0.45
36:5:996:A:C2	36:5:997:A:H1'	2.52	0.45
1:6:1292:G:C5	1:6:1293:U:C4	3.05	0.45
1:6:246:G:C6	1:6:247:A:C6	3.04	0.45
1:6:567:A:N1	1:6:583:C:H1'	2.31	0.45
13:C1:13:PHE:CD2	13:C1:15:LYS:HB3	2.52	0.45
21:C9:28:LEU:HD12	21:C9:29:GLU:H	1.82	0.45
26:D4:23:PHE:HE2	26:D4:75:VAL:HG12	1.81	0.45
36:1:860:G:C6	39:L2:181:LYS:HB2	2.52	0.45
40:L3:139:GLN:H	40:L3:139:GLN:HG3	2.99	0.45
41:L4:99:MET:SD	41:L4:102:PRO:HA	2.57	0.45
41:L4:138:ARG:HE	41:L4:240:PRO:HD2	1.92	0.45
41:L4:233:LEU:HD13	41:L4:238:LEU:HD11	3.31	0.45
43:L6:52:VAL:HG22	43:L6:67:GLY:HA2	2.27	0.45
43:L6:65:ILE:HA	43:L6:65:ILE:HD13	3.95	0.45
46:L9:44:THR:HG21	50:M4:12:TRP:CH2	4.06	0.45
47:M0:19:LYS:HB2	47:M0:26:VAL:HG11	3.57	0.45
51:M5:18:VAL:O	51:M5:22:LEU:HD22	2.16	0.45
54:M8:81:VAL:HG22	54:M8:101:VAL:HG22	1.98	0.45
61:N5:82:LEU:HB3	61:N5:84:PHE:CE2	2.52	0.45
62:N6:30:LEU:C	62:N6:32:SER:H	3.09	0.45
63:N7:53:VAL:HG21	63:N7:62:VAL:HG13	1.99	0.45
67:O1:70:ARG:O	67:O1:71:LEU:HD23	2.41	0.45
69:O3:13:HIS:NE2	69:O3:28:SER:OG	2.56	0.45
71:O5:31:LEU:HD22	71:O5:41:LEU:HD21	2.33	0.45
71:O5:31:LEU:O	71:O5:34:GLN:HB2	2.17	0.45
72:O6:34:SER:OG	72:O6:37:THR:OG1	2.23	0.45
72:O6:91:ASN:O	72:O6:94:ILE:HG22	4.60	0.45
3:S1:156:ALA:HB1	3:S1:160:HIS:HB2	2.92	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:S4:49:ARG:HG3	6:S4:55:ALA:HB3	1.98	0.45
9:S7:133:THR:HG22	9:S7:157:LYS:O	3.95	0.45
10:S8:170:SER:OG	10:S8:181:GLY:HA2	2.16	0.45
11:S9:53:ARG:O	11:S9:57:ARG:N	2.90	0.45
34:SR:38:ARG:HG2	34:SR:67:ILE:CG2	2.47	0.45
36:1:1275:C:H2'	36:1:1276:U:O4'	2.16	0.45
36:1:1911:A:H8	36:1:1911:A:O5'	2.00	0.45
36:1:2295:A:OP1	59:N3:63:LYS:NZ	2.45	0.45
36:1:874:U:OP1	40:L3:241:LYS:HG2	2.17	0.45
1:2:1086:A:C6	1:2:1087:A:C6	3.05	0.45
1:2:1165:G:O6	1:2:1166:A:N6	2.50	0.45
36:5:1571:A:H1'	36:5:1572:U:C5	2.52	0.45
36:5:1781:C:H2'	36:5:1782:U:H6	1.81	0.45
36:5:2931:C:H2'	36:5:2932:U:O4'	2.17	0.45
44:L7:60:ARG:NH2	36:5:516:A:O3'	303.00	0.45
1:6:1765:A:H5'	1:6:1767:G:N7	2.32	0.45
13:C1:10:GLU:HG2	1:6:327:U:H1'	270.13	0.45
1:6:53:G:H2'	1:6:54:C:C6	2.52	0.45
1:6:547:U:H2'	1:6:548:G:O4'	2.17	0.45
37:7:11:A:H4'	37:7:13:A:C8	2.51	0.45
16:C4:19:ILE:HG22	16:C4:21:ALA:HB2	1.99	0.45
5:S3:211:PRO:HG3	19:C7:19:ARG:HB3	5.23	0.45
19:C7:66:VAL:C	19:C7:68:GLY:H	4.35	0.45
19:C7:84:TYR:O	19:C7:86:PRO:HD3	2.16	0.45
22:D0:72:ASN:OD1	22:D0:73:GLY:N	2.64	0.45
2:S0:7:PHE:HE1	23:D1:39:VAL:HG21	4.62	0.45
23:D1:74:GLN:HG3	23:D1:79:LEU:O	2.17	0.45
26:D4:123:LYS:H	26:D4:123:LYS:HG2	1.54	0.45
43:L6:137:ASP:O	43:L6:141:VAL:HG23	2.17	0.45
44:L7:92:ILE:HA	44:L7:92:ILE:HD12	1.76	0.45
45:L8:145:ASN:C	45:L8:147:LYS:H	3.55	0.45
46:L9:161:LEU:HD22	46:L9:179:ILE:HD12	1.99	0.45
47:M0:205:SER:HB3	47:M0:208:ASN:OD1	2.17	0.45
47:M0:36:LEU:HD13	47:M0:87:LEU:HD13	1.99	0.45
48:M1:101:ASN:H	48:M1:101:ASN:ND2	2.13	0.45
53:M7:136:ILE:O	53:M7:137:ASN:ND2	2.90	0.45
53:M7:30:ARG:CZ	53:M7:34:GLN:HG3	2.47	0.45
56:N0:137:ARG:HG2	56:N0:139:TYR:CE1	2.52	0.45
60:N4:82:ILE:O	60:N4:85:ALA:N	4.39	0.45
63:N7:10:VAL:HG22	63:N7:24:VAL:HG13	1.98	0.45
63:N7:42:LEU:H	63:N7:42:LEU:HD12	2.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:965:A:C2	64:N8:43:ILE:HD12	2.47	0.45
68:O2:100:ILE:O	68:O2:105:ARG:NH1	3.10	0.45
68:O2:15:LYS:HB3	68:O2:15:LYS:HE3	3.07	0.45
71:O5:49:LYS:HA	71:O5:49:LYS:HD3	1.76	0.45
78:Q2:93:LEU:H	78:Q2:93:LEU:HD23	1.82	0.45
11:S9:114:TYR:C	11:S9:116:LEU:H	2.96	0.45
11:S9:6:ARG:HD2	1:6:771:A:O2'	389.68	0.45
35:SM:27:LYS:HB2	48:M1:46:VAL:HG21	2.63	0.45
34:SR:85:TRP:CD1	34:SR:109:ASP:HB3	3.51	0.45
36:1:1069:C:H2'	36:1:1070:U:C6	2.53	0.44
36:1:1366:A:C2	36:1:1367:G:C4	3.06	0.44
36:1:1471:U:H2'	36:1:1472:U:H6	1.82	0.44
36:1:2278:C:H2'	36:1:2279:A:H5''	1.99	0.44
36:1:2780:A:H2'	36:1:2781:U:C6	2.52	0.44
36:1:314:U:H2'	36:1:315:C:C6	2.53	0.44
36:1:729:C:H2'	36:1:730:C:H6	1.83	0.44
1:2:1066:C:O3'	3:S1:149:GLN:HG3	2.17	0.44
1:2:1173:C:H2'	1:2:1174:C:C6	2.52	0.44
1:2:1393:C:H2'	1:2:1394:G:O4'	2.17	0.44
1:2:1470:C:OP1	1:2:1540:G:O2'	2.35	0.44
1:2:374:U:OP1	13:C1:96:LYS:NZ	2.36	0.44
1:2:890:C:H2'	1:2:891:A:H8	1.82	0.44
38:4:23:U:H4'	62:N6:17:LYS:HG2	1.99	0.44
36:5:1190:A:C5	36:5:1193:A:H1'	2.52	0.44
36:5:1579:C:C5'	36:5:1649:U:H5''	2.47	0.44
36:5:2172:A:C6	36:5:2173:U:C4	3.05	0.44
36:5:240:U:O2'	36:5:241:G:H8	2.00	0.44
36:5:2448:G:H22	36:5:2498:U:H3	1.64	0.44
41:L4:82:THR:OG1	36:5:365:A:H1'	121.81	0.44
1:6:151:G:N2	1:6:164:A:C4	2.85	0.44
1:6:1589:C:H2'	1:6:1590:G:C8	2.52	0.44
1:6:1779:U:H2'	1:6:1781:A:OP2	2.17	0.44
12:C0:32:HIS:CD2	12:C0:33:GLU:H	2.73	0.44
13:C1:46:LYS:HG3	13:C1:50:GLU:CD	5.74	0.44
14:C2:61:VAL:HA	14:C2:89:ILE:HG22	1.99	0.44
21:C9:117:SER:HB3	21:C9:123:ARG:HB3	2.34	0.44
21:C9:66:TYR:HA	21:C9:124:ILE:HB	1.99	0.44
22:D0:26:LEU:HB2	22:D0:89:ARG:HB2	2.92	0.44
25:D3:69:ARG:HD2	25:D3:117:ILE:HG12	5.25	0.44
26:D4:78:SER:OG	26:D4:79:VAL:N	2.51	0.44
39:L2:191:LEU:HD23	36:5:1794:G:H4'	192.20	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:L2:96:LEU:O	79:Q3:87:ARG:HD3	2.17	0.44
40:L3:87:VAL:HB	40:L3:110:LEU:HD11	1.99	0.44
41:L4:180:LYS:HA	36:5:1386:A:N3	118.74	0.44
41:L4:271:LYS:O	41:L4:273:GLY:N	2.68	0.44
43:L6:65:ILE:O	43:L6:76:LEU:HA	2.58	0.44
44:L7:203:TRP:CD1	44:L7:204:PRO:HD2	2.52	0.44
44:L7:90:LYS:HZ2	44:L7:91:GLY:N	4.40	0.44
45:L8:97:TYR:HB3	45:L8:131:ALA:HA	1.98	0.44
45:L8:97:TYR:OH	45:L8:204:ARG:N	2.47	0.44
36:1:3198:U:H1'	46:L9:21:LYS:HB2	2.00	0.44
48:M1:166:LYS:C	48:M1:168:ASP:H	2.31	0.44
49:M3:58:VAL:CG1	36:5:75:G:H5''	87.93	0.44
52:M6:58:LEU:HD12	52:M6:58:LEU:HA	1.83	0.44
53:M7:108:ASP:O	53:M7:110:THR:N	2.47	0.44
41:L4:282:SER:CB	54:M8:126:GLN:HE21	2.47	0.44
55:M9:170:ARG:O	55:M9:174:ALA:N	2.47	0.44
60:N4:82:ILE:O	60:N4:84:GLY:N	3.82	0.44
61:N5:67:ILE:HD12	61:N5:121:LYS:HG3	1.99	0.44
71:O5:102:GLU:HG3	71:O5:106:LYS:HE3	4.90	0.44
46:L9:176:LEU:HD13	76:Q0:86:ALA:HB3	2.28	0.44
76:Q0:92:ASP:O	76:Q0:105:PRO:HG3	2.17	0.44
1:2:1116:A:P	77:Q1:17:ARG:HH21	2.39	0.44
79:Q3:6:LYS:HZ2	79:Q3:7:LYS:HG3	7.86	0.44
2:S0:74:VAL:HG22	2:S0:96:THR:HG23	2.62	0.44
3:S1:30:PHE:CD1	3:S1:96:LEU:HD22	2.52	0.44
4:S2:144:TRP:CZ2	4:S2:173:PRO:HG3	2.95	0.44
6:S4:182:TYR:CZ	6:S4:192:ILE:HD11	4.11	0.44
11:S9:78:ARG:O	11:S9:82:ARG:HB2	3.21	0.44
34:SR:242:SER:O	34:SR:292:LEU:HD21	2.17	0.44
36:1:1325:U:H2'	36:1:1326:A:C8	2.52	0.44
36:1:2273:G:H22	36:1:2311:G:H2'	1.82	0.44
36:1:2381:G:C2'	36:1:2382:G:H5'	2.47	0.44
36:1:3356:G:H2'	36:1:3357:U:O4'	2.17	0.44
86:1:3401:SPS:H91	86:1:3401:SPS:O1	2.18	0.44
36:1:564:G:H2'	36:1:565:U:C6	2.52	0.44
1:2:116:U:H2'	1:2:117:U:C6	2.53	0.44
1:2:1173:C:H2'	1:2:1174:C:H6	1.82	0.44
1:2:1408:G:H2'	1:2:1409:G:O4'	2.17	0.44
1:2:144:U:C2	1:2:145:A:C8	3.05	0.44
1:2:350:U:H5''	1:2:352:A:H5'	1.98	0.44
1:2:413:U:H2'	1:2:414:C:C6	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:648:G:C2	1:2:687:G:N2	2.85	0.44
1:2:700:C:H42	1:2:738:G:H1	1.65	0.44
74:O8:17:ARG:NH2	36:5:1824:U:H4'	140.86	0.44
36:5:2697:A:H2'	36:5:2698:G:C8	2.51	0.44
36:5:3041:U:H2'	36:5:3042:U:C6	2.52	0.44
73:O7:13:ASN:HD21	36:5:901:G:H5''	145.07	0.44
64:N8:13:GLY:CA	36:5:943:U:H3'	163.73	0.44
1:6:486:G:H22	1:6:501:U:H3	1.66	0.44
1:6:538:A:H2	1:6:540:G:N2	2.15	0.44
1:6:53:G:C2	1:6:54:C:C2	3.05	0.44
38:8:81:U:OP1	38:8:81:U:H4'	2.16	0.44
13:C1:35:TYR:HD1	13:C1:49:ILE:HG12	3.71	0.44
14:C2:54:ARG:HB3	33:E1:128:ALA:HB1	7.66	0.44
16:C4:84:ARG:HB2	16:C4:118:VAL:HG23	1.98	0.44
1:2:927:C:H4'	16:C4:125:SER:OG	2.18	0.44
20:C8:57:ARG:HB2	20:C8:60:GLU:HG3	1.99	0.44
24:D2:93:LEU:O	24:D2:94:LEU:HD23	2.74	0.44
29:D7:28:PRO:C	29:D7:30:SER:H	2.78	0.44
29:D7:2:VAL:C	29:D7:4:VAL:H	2.21	0.44
30:D8:32:PHE:CE2	30:D8:38:ARG:HD2	2.52	0.44
33:E1:93:HIS:CG	33:E1:94:LYS:H	2.34	0.44
39:L2:27:ALA:HB3	39:L2:128:ARG:NH2	3.40	0.44
40:L3:147:GLU:O	40:L3:151:ILE:HG13	2.89	0.44
41:L4:141:ARG:C	41:L4:143:GLU:N	4.05	0.44
41:L4:140:HIS:HA	41:L4:177:ASP:OD1	2.17	0.44
41:L4:307:GLN:OE1	36:5:1346:G:O2'	202.15	0.44
44:L7:239:LEU:O	44:L7:242:SER:OG	2.29	0.44
47:M0:3:ARG:HA	47:M0:3:ARG:HD2	3.12	0.44
48:M1:36:VAL:HG21	48:M1:123:PHE:HD2	1.83	0.44
50:M4:24:LYS:HB3	50:M4:24:LYS:HE2	4.57	0.44
52:M6:31:GLN:HE21	52:M6:32:LYS:H	1.65	0.44
59:N3:27:ASP:O	59:N3:29:SER:N	3.95	0.44
60:N4:4:GLU:HG2	60:N4:30:ARG:CD	2.47	0.44
61:N5:64:GLU:O	61:N5:65:GLN:HG3	2.17	0.44
63:N7:29:HIS:CE1	63:N7:42:LEU:HD13	2.52	0.44
66:O0:13:LYS:O	66:O0:17:VAL:HG23	2.17	0.44
68:O2:69:SER:HG	68:O2:71:HIS:CG	2.35	0.44
71:O5:118:ILE:O	71:O5:119:LYS:HD2	5.07	0.44
73:O7:18:LEU:HD21	75:O9:51:ILE:HG22	1.99	0.44
75:O9:24:PRO:HB2	75:O9:27:ILE:HD13	2.00	0.44
36:1:2150:G:H4'	79:Q3:22:LEU:HD21	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S2:187:LEU:HD21	4:S2:215:PHE:HB2	1.99	0.44
4:S2:67:GLN:OE1	4:S2:67:GLN:N	2.85	0.44
6:S4:136:VAL:HG11	6:S4:148:ARG:NH2	2.31	0.44
9:S7:117:THR:HG23	1:6:639:U:P	364.02	0.44
35:SM:61:ILE:HG13	35:SM:61:ILE:H	1.61	0.44
34:SR:117:LYS:H	34:SR:117:LYS:HD2	1.82	0.44
36:1:1282:G:H2'	36:1:1283:C:O4'	2.17	0.44
36:1:1369:A:H2'	36:1:1370:G:O4'	2.18	0.44
36:1:1764:U:H3'	36:1:1765:U:O4'	2.17	0.44
36:1:2273:G:N2	36:1:2311:G:H2'	2.32	0.44
36:1:957:C:H4'	36:1:2799:A:C5	2.52	0.44
36:1:307:A:H2'	36:1:308:A:H8	1.83	0.44
1:2:1584:G:N2	1:2:1611:A:OP2	2.20	0.44
1:2:1698:G:H1'	1:2:1699:G:OP1	2.17	0.44
1:2:1701:A:H3'	1:2:1702:A:H5''	1.99	0.44
1:2:1729:C:H2'	1:2:1730:A:O4'	2.17	0.44
1:2:432:G:H2'	1:2:433:C:O4'	2.18	0.44
1:2:628:G:N1	1:2:970:A:OP2	2.37	0.44
38:4:128:U:OP1	38:4:129:C:N4	2.38	0.44
36:5:1574:C:O5'	36:5:1574:C:H6	2.00	0.44
36:5:1587:A:O2'	36:5:1590:G:O6	2.31	0.44
36:5:2133:U:O4	36:5:2147:A:H2	1.99	0.44
36:5:2442:G:H1	36:5:2505:U:H3	1.66	0.44
36:5:26:A:N3	36:5:328:U:O2'	2.42	0.44
36:5:2771:U:N3	36:5:2773:C:N3	2.65	0.44
36:5:348:A:H1'	36:5:352:A:O2'	2.18	0.44
1:6:1111:G:C2	1:6:1112:G:H1'	2.52	0.44
1:6:1244:A:N3	1:6:1244:A:H3'	2.32	0.44
1:6:1528:U:H2'	1:6:1529:C:O4'	2.17	0.44
1:6:296:U:C4	1:6:297:U:C4	3.06	0.44
1:6:912:U:H4'	1:6:913:G:H5'	1.99	0.44
73:O7:65:ARG:HH22	38:8:43:A:N6	84.80	0.44
20:C8:91:ASP:CG	20:C8:94:ASP:HB3	5.29	0.44
22:D0:27:THR:HG23	22:D0:113:ASP:HB3	1.99	0.44
26:D4:61:ARG:HG3	26:D4:62:THR:N	3.44	0.44
28:D6:37:LYS:HD3	28:D6:72:HIS:CD2	2.53	0.44
28:D6:51:ARG:NH2	28:D6:55:GLU:OE1	2.65	0.44
3:S1:111:ARG:HB3	28:D6:68:TYR:CG	3.74	0.44
42:L5:260:PHE:HB3	42:L5:264:GLN:HB3	2.36	0.44
43:L6:44:ALA:O	43:L6:48:ARG:HB3	2.91	0.44
46:L9:87:LYS:HB2	46:L9:187:ILE:HG12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:L9:187:ILE:HG22	46:L9:188:THR:H	1.81	0.44
50:M4:15:VAL:HG13	56:N0:150:PHE:O	2.34	0.44
51:M5:59:PHE:CD1	51:M5:133:ILE:HD11	2.47	0.44
51:M5:178:HIS:CE1	51:M5:179:LYS:HG3	2.53	0.44
53:M7:13:LYS:HE2	53:M7:152:GLU:HB2	3.01	0.44
55:M9:8:LYS:HE3	55:M9:22:VAL:HG23	1.99	0.44
56:N0:40:ARG:HG2	56:N0:40:ARG:HH11	4.28	0.44
56:N0:46:GLN:O	37:7:77:G:H5''	301.20	0.44
57:N1:11:THR:CG2	57:N1:55:LYS:HD2	2.54	0.44
64:N8:65:GLN:O	64:N8:66:ALA:HB3	3.08	0.44
65:N9:22:LYS:HG2	65:N9:22:LYS:H	1.54	0.44
66:O0:24:THR:OG1	66:O0:29:SER:OG	3.30	0.44
66:O0:92:ILE:HG21	66:O0:100:ILE:HD11	3.79	0.44
68:O2:111:ARG:NE	68:O2:115:LEU:HD11	2.33	0.44
69:O3:45:LEU:HD23	69:O3:71:VAL:HG12	2.00	0.44
70:O4:20:ILE:HD11	70:O4:32:ALA:HB1	1.99	0.44
73:O7:45:ARG:HH11	73:O7:47:TYR:HE2	1.66	0.44
48:M1:62:ASN:ND2	78:Q2:101:GLY:O	2.42	0.44
79:Q3:49:ARG:HD3	79:Q3:51:ALA:O	4.44	0.44
2:S0:148:ASP:HB2	2:S0:164:ASN:ND2	2.33	0.44
2:S0:63:ILE:O	2:S0:67:ILE:HD13	3.38	0.44
3:S1:124:ASN:HB3	3:S1:138:PHE:HA	2.00	0.44
3:S1:191:GLU:HB2	3:S1:194:ASN:CG	2.37	0.44
5:S3:125:TYR:O	5:S3:129:SER:OG	2.93	0.44
6:S4:107:GLY:HA2	6:S4:189:LEU:HG	2.50	0.44
6:S4:184:THR:C	6:S4:189:LEU:HD13	3.70	0.44
7:S5:201:ALA:HA	7:S5:211:ILE:HD11	3.60	0.44
9:S7:129:LEU:HA	9:S7:129:LEU:HD23	1.75	0.44
11:S9:102:GLU:HA	11:S9:105:LEU:HB2	2.50	0.44
34:SR:221:MET:HA	34:SR:233:THR:HG23	3.18	0.44
36:1:1094:U:H1'	36:1:1096:U:H2'	2.00	0.44
36:1:3084:C:OP1	60:N4:38:SER:OG	2.25	0.44
36:1:3197:G:H2'	36:1:3198:U:H5''	1.98	0.44
36:1:3290:G:C6	36:1:3291:G:C5	3.05	0.44
36:1:643:U:O4	36:1:644:G:C6	2.70	0.44
36:1:954:U:O4	36:1:1115:G:H1'	2.18	0.44
1:2:1352:G:H2'	1:2:1353:U:O4'	2.17	0.44
1:2:1555:A:OP2	17:C5:47:ARG:NH2	2.50	0.44
1:2:47:A:N1	1:2:386:G:H1'	2.33	0.44
1:2:811:A:C2	1:2:858:G:H1'	2.52	0.44
36:5:1093:A:N3	36:5:1096:U:N3	2.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:1549:U:H2'	36:5:1550:C:C6	2.52	0.44
36:5:184:U:H2'	36:5:185:C:C6	2.52	0.44
36:5:2628:A:H1'	36:5:2798:C:C2	2.53	0.44
36:5:3236:U:H2'	36:5:3237:U:C6	2.53	0.44
1:6:119:A:H1'	1:6:397:A:C4	2.52	0.44
1:6:1484:G:H2'	1:6:1485:C:H6	1.82	0.44
1:6:151:G:H22	1:6:163:G:N2	2.14	0.44
1:6:1607:G:H2'	1:6:1608:U:H6	1.83	0.44
1:6:1742:U:H2'	1:6:1743:U:H6	1.82	0.44
1:6:1754:A:H4'	1:6:1755:A:O5'	2.16	0.44
1:6:291:G:H2'	1:6:292:U:C6	2.52	0.44
11:S9:133:HIS:NE2	1:6:512:A:O3'	447.10	0.44
26:D4:11:LYS:HD3	1:6:784:C:N4	417.15	0.44
38:8:78:G:H2'	38:8:79:A:O4'	2.17	0.44
13:C1:72:THR:HG22	13:C1:124:THR:HA	2.57	0.44
10:S8:10:LYS:HG2	13:C1:133:LYS:CE	2.66	0.44
18:C6:109:PHE:O	18:C6:113:ASP:N	2.47	0.44
1:2:1586:A:H5''	18:C6:136:SER:HB2	1.99	0.44
25:D3:41:SER:HA	25:D3:42:PRO:HD3	1.77	0.44
25:D3:91:GLY:O	25:D3:93:LEU:N	2.47	0.44
25:D3:96:VAL:HG23	25:D3:97:ASP:H	1.82	0.44
28:D6:44:ILE:HD12	28:D6:45:VAL:HG13	2.00	0.44
32:E0:31:LYS:HG2	32:E0:31:LYS:H	1.59	0.44
39:L2:109:GLU:H	39:L2:109:GLU:CD	4.23	0.44
39:L2:147:ARG:HH12	39:L2:155:LYS:HD2	1.81	0.44
39:L2:218:HIS:HD2	36:5:2246:G:OP1	221.68	0.44
40:L3:183:LEU:O	40:L3:191:LYS:NZ	2.84	0.44
40:L3:21:ARG:HG2	40:L3:269:GLN:HG2	1.99	0.44
40:L3:77:THR:HG23	40:L3:327:CYS:HA	1.98	0.44
40:L3:67:PHE:CE1	59:N3:88:ARG:HB3	2.52	0.44
37:3:44:C:H4'	42:L5:152:ARG:HG3	1.97	0.44
45:L8:143:ILE:HB	45:L8:169:LEU:HD13	2.58	0.44
48:M1:45:PRO:HB2	48:M1:67:VAL:HG22	6.24	0.44
49:M3:50:PRO:HB3	49:M3:140:SER:O	2.17	0.44
51:M5:94:TYR:CE2	51:M5:96:ARG:HB2	3.72	0.44
53:M7:13:LYS:HB3	53:M7:13:LYS:HE2	2.55	0.44
55:M9:117:LYS:HA	55:M9:120:TYR:HB3	1.99	0.44
57:N1:84:TYR:HD1	65:N9:23:LYS:HA	3.96	0.44
59:N3:48:ARG:HG2	36:5:2339:C:P	245.81	0.44
61:N5:72:ALA:O	61:N5:76:VAL:HG23	2.17	0.44
64:N8:47:LYS:HE2	64:N8:48:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
69:O3:71:VAL:HG13	69:O3:81:VAL:CG1	2.48	0.44
70:O4:37:LYS:HB2	70:O4:58:ARG:NH2	2.33	0.44
72:O6:15:LYS:NZ	72:O6:16:LYS:H	5.91	0.44
36:1:2896:A:H4'	76:Q0:95:VAL:HG11	1.99	0.44
78:Q2:12:CYS:SG	78:Q2:74:CYS:SG	3.28	0.44
2:S0:80:THR:HA	2:S0:83:GLN:HE21	6.05	0.44
3:S1:157:GLN:HB2	3:S1:160:HIS:NE2	4.11	0.44
4:S2:66:PHE:HB3	4:S2:130:ILE:HG23	2.99	0.44
5:S3:70:THR:HG22	5:S3:86:LEU:HD13	1.98	0.44
6:S4:88:ASP:HA	6:S4:122:LYS:HE2	1.99	0.44
7:S5:25:LEU:HD22	7:S5:25:LEU:H	1.82	0.44
8:S6:20:ASP:O	8:S6:24:ILE:HG13	2.30	0.44
9:S7:6:ALA:HB1	9:S7:9:LEU:HD12	1.97	0.44
11:S9:129:ILE:HA	11:S9:134:ILE:CG1	3.81	0.44
11:S9:53:ARG:O	11:S9:57:ARG:HG2	2.17	0.44
36:1:2590:A:C4	36:1:2591:A:C8	3.06	0.44
36:1:2709:C:H2'	36:1:2710:C:C6	2.52	0.44
36:1:531:G:H2'	36:1:532:A:C8	2.52	0.44
36:1:874:U:H5''	36:1:2950:G:OP1	2.18	0.44
36:1:869:G:H1'	36:1:891:G:N2	2.32	0.44
36:1:929:A:H2'	36:1:930:U:C6	2.53	0.44
1:2:1434:U:H4'	31:D9:24:CYS:HB2	2.00	0.44
1:2:1450:U:H2'	1:2:1451:C:C6	2.53	0.44
1:2:1460:A:OP2	35:SM:68:ARG:HD3	2.18	0.44
1:2:223:U:H2'	1:2:224:C:C6	2.52	0.44
1:2:526:A:H2'	1:2:527:A:O4'	2.17	0.44
1:2:824:G:N2	1:2:849:C:O2	2.51	0.44
36:5:1626:U:O3'	36:5:1632:A:H4'	2.17	0.44
36:5:1915:A:H2'	36:5:1916:U:C6	2.51	0.44
36:5:209:A:H4'	36:5:211:A:N7	2.32	0.44
40:L3:259:HIS:HB3	36:5:2987:A:O2'	216.63	0.44
36:5:3133:C:H2'	36:5:3134:A:O4'	2.17	0.44
65:N9:14:ARG:NH2	36:5:952:A:OP1	208.07	0.44
1:6:1470:C:H42	1:6:1574:G:H1	1.66	0.44
1:6:446:A:N6	1:6:461:G:H21	2.15	0.44
1:6:481:A:H2'	1:6:482:U:O4'	2.17	0.44
1:6:560:U:H2'	1:6:561:G:H8	1.83	0.44
11:S9:79:ARG:NH1	1:6:762:A:OP1	407.55	0.44
1:6:842:C:H2'	1:6:843:U:O4'	2.17	0.44
75:O9:21:ARG:NH2	38:8:52:A:OP1	79.45	0.44
14:C2:61:VAL:HG13	14:C2:121:VAL:HB	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:C3:2:GLY:O	15:C3:3:ARG:HB3	2.18	0.44
17:C5:26:LEU:HD23	17:C5:87:PRO:HB2	8.18	0.44
20:C8:142:GLY:O	20:C8:145:ARG:HD3	2.17	0.44
25:D3:133:LEU:HD21	25:D3:137:LYS:HE3	1.98	0.44
26:D4:21:LYS:NZ	26:D4:77:ASN:OD1	2.49	0.44
7:S5:112:ARG:HH22	27:D5:94:LYS:HD3	1.83	0.44
1:2:567:A:H4'	32:E0:10:ARG:O	2.17	0.44
33:E1:86:THR:HG23	33:E1:87:THR:H	4.26	0.44
39:L2:36:GLU:HG2	39:L2:90:ALA:O	2.18	0.44
42:L5:184:ASP:OD1	42:L5:187:THR:N	2.50	0.44
42:L5:277:LEU:HB3	42:L5:281:GLU:OE2	5.54	0.44
36:1:503:C:OP1	43:L6:26:ARG:NH1	2.51	0.44
44:L7:80:GLN:OE1	57:N1:136:ARG:N	2.41	0.44
45:L8:158:ASP:HA	45:L8:159:PRO:C	4.55	0.44
36:1:2645:G:OP2	47:M0:117:GLY:HA2	2.18	0.44
45:L8:137:ASN:HB3	51:M5:2:GLY:HA2	3.96	0.44
52:M6:111:PRO:O	52:M6:113:ASP:N	2.50	0.44
52:M6:14:HIS:CD2	52:M6:124:LEU:HD13	2.52	0.44
54:M8:165:ILE:HD13	54:M8:166:LEU:N	5.30	0.44
55:M9:154:ALA:O	55:M9:158:GLU:HG2	2.59	0.44
44:L7:77:VAL:HG22	57:N1:139:ARG:HB2	4.28	0.44
47:M0:168:SER:HB2	57:N1:160:ILE:C	4.29	0.44
58:N2:24:GLU:HG3	58:N2:25:ASN:OD1	2.17	0.44
59:N3:26:ALA:HB3	59:N3:101:VAL:HG13	1.99	0.44
67:O1:51:LEU:HB3	67:O1:55:LEU:HD12	2.61	0.44
79:Q3:6:LYS:HG2	79:Q3:7:LYS:HG3	4.77	0.44
3:S1:32:ILE:O	3:S1:43:VAL:HB	2.18	0.44
4:S2:228:ASN:N	4:S2:228:ASN:OD1	2.67	0.44
6:S4:139:VAL:HG13	6:S4:150:PRO:HG3	1.99	0.44
6:S4:29:PRO:HG3	6:S4:45:ILE:HD13	1.99	0.44
7:S5:153:GLY:O	7:S5:155:ALA:N	2.96	0.44
11:S9:142:ASN:O	11:S9:143:ILE:HD12	3.71	0.44
36:1:2100:A:H5'	36:1:2101:C:OP1	2.18	0.44
36:1:2134:G:H1	36:1:2146:C:H42	1.64	0.44
36:1:420:G:N2	36:1:2385:G:OP2	2.44	0.44
36:1:2510:U:O2'	36:1:2511:A:H8	2.01	0.44
36:1:295:A:H2'	36:1:296:A:O4'	2.17	0.44
36:1:426:G:H5'	68:O2:50:ILE:HG22	2.00	0.44
36:1:428:A:H2'	36:1:429:U:C6	2.52	0.44
36:1:543:C:N4	36:1:548:G:H1	2.16	0.44
36:1:75:G:H5'	49:M3:59:ARG:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1299:G:H2'	1:2:1300:A:C8	2.52	0.44
1:2:1582:U:OP1	18:C6:135:ARG:NH1	2.50	0.44
1:2:409:C:H2'	1:2:410:A:H8	1.83	0.44
1:2:737:A:OP2	1:2:737:A:H2'	2.17	0.44
1:2:766:U:H5'	1:2:767:U:H5''	2.00	0.44
38:4:77:A:H2'	38:4:78:G:O4'	2.18	0.44
36:5:1288:U:H2'	36:5:1289:G:H8	1.83	0.44
36:5:1313:G:H2'	36:5:1314:C:C6	2.52	0.44
36:5:1440:G:H2'	36:5:1441:G:H8	1.83	0.44
36:5:1711:C:H2'	36:5:1712:G:O4'	2.18	0.44
36:5:326:U:H6	36:5:326:U:O5'	1.99	0.44
36:5:662:U:H2'	36:5:663:C:C6	2.52	0.44
36:5:746:A:H2'	36:5:747:A:C8	2.52	0.44
1:6:123:G:H2'	1:6:124:A:O4'	2.18	0.44
1:6:257:A:H2'	1:6:258:C:C6	2.53	0.44
1:6:763:G:C6	1:6:764:U:C4	3.06	0.44
42:L5:265:TYR:OH	37:7:121:U:OP2	312.89	0.44
38:8:63:G:OP1	38:8:90:U:H5'	2.18	0.44
12:C0:11:ILE:HD12	12:C0:42:VAL:HG13	1.99	0.44
13:C1:57:LYS:HD3	13:C1:131:ILE:HG23	3.68	0.44
15:C3:26:PHE:CE2	15:C3:28:LEU:HB2	2.76	0.44
27:D5:40:VAL:HA	27:D5:75:LEU:HD11	1.99	0.44
27:D5:93:SER:HB3	27:D5:100:ILE:HB	1.98	0.44
27:D5:96:SER:C	27:D5:98:GLN:H	2.21	0.44
30:D8:41:VAL:O	30:D8:62:GLU:HB2	2.18	0.44
39:L2:214:GLY:O	39:L2:215:ASN:HB2	4.68	0.44
41:L4:35:VAL:HG11	41:L4:244:LEU:HD21	2.35	0.44
42:L5:43:LYS:HB3	42:L5:46:THR:OG1	3.65	0.44
42:L5:60:ILE:HB	42:L5:80:SER:HB3	3.51	0.44
43:L6:62:THR:HG21	43:L6:78:ARG:HB3	3.20	0.44
44:L7:96:PRO:O	44:L7:100:ARG:HB2	2.27	0.44
45:L8:163:VAL:HG22	45:L8:166:LEU:HD12	2.34	0.44
45:L8:81:THR:HB	45:L8:181:LYS:HB2	3.77	0.44
48:M1:92:ARG:HB2	48:M1:95:ASN:ND2	2.32	0.44
49:M3:76:THR:HG22	49:M3:101:ARG:HB3	2.00	0.44
49:M3:157:ARG:HG2	49:M3:158:ALA:N	2.33	0.44
53:M7:108:ASP:C	53:M7:110:THR:H	2.20	0.44
41:L4:302:ALA:HB2	54:M8:39:ARG:HH22	2.85	0.44
55:M9:173:ARG:O	55:M9:177:VAL:HG23	2.19	0.44
62:N6:124:GLY:C	62:N6:126:LEU:H	3.55	0.44
62:N6:52:ARG:O	62:N6:53:ASP:HB2	4.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:90:GLU:HA	63:N7:93:LYS:HB2	1.98	0.44
68:O2:60:ASN:O	68:O2:64:LYS:HB2	2.17	0.44
61:N5:117:ASN:HA	75:O9:14:ALA:HB1	2.00	0.44
2:S0:88:LYS:HG3	2:S0:201:LEU:O	4.75	0.44
4:S2:152:HIS:CG	4:S2:174:ARG:HG2	2.52	0.44
2:S0:119:ARG:NH1	4:S2:241:ASP:OD2	2.50	0.44
6:S4:198:LYS:HB3	6:S4:198:LYS:HE2	3.36	0.44
6:S4:47:PHE:HD2	6:S4:48:LEU:HD12	1.82	0.44
8:S6:67:VAL:O	8:S6:68:LEU:HB2	2.46	0.44
11:S9:40:LYS:HB2	11:S9:40:LYS:HE3	1.65	0.44
36:1:109:A:H4'	36:1:110:G:OP1	2.17	0.44
36:1:1364:C:H5''	54:M8:3:ILE:HD13	1.99	0.44
36:1:1631:C:H5''	36:1:1632:A:H5''	2.00	0.44
36:1:1694:U:O2'	36:1:1695:U:H5'	2.18	0.44
36:1:2190:U:C4	36:1:2191:U:C4	3.06	0.44
36:1:2413:A:H2'	36:1:2414:G:C8	2.53	0.44
36:1:2502:A:N3	36:1:2502:A:H2'	2.33	0.44
36:1:2656:A:C4	36:1:2658:G:N7	2.86	0.44
36:1:2759:U:H6	36:1:2759:U:O5'	2.01	0.44
36:1:2818:U:C6	36:1:2818:U:H5'	2.44	0.44
36:1:352:A:N6	36:1:365:A:H5''	2.29	0.44
1:2:1132:A:H2'	1:2:1133:A:H8	1.81	0.44
1:2:1642:G:H2'	1:2:1643:U:H6	1.83	0.44
37:3:42:A:H2'	37:3:43:U:H6	1.83	0.44
38:4:149:A:N3	45:L8:55:TYR:OH	2.39	0.44
36:5:1062:A:H5''	36:5:1063:G:H5'	1.99	0.44
68:O2:99:ASN:HB2	36:5:1388:U:O2'	131.17	0.44
36:5:1477:A:OP1	36:5:3075:G:O2'	2.33	0.44
36:5:1733:G:H2'	36:5:1734:G:C8	2.52	0.44
36:5:2315:G:C2	36:5:2316:G:N7	2.86	0.44
36:5:249:U:OP2	36:5:249:U:H6	2.01	0.44
36:5:2781:U:C4	36:5:2782:U:C4	3.05	0.44
36:5:2941:A:O5'	36:5:2943:G:H4'	2.17	0.44
36:5:437:G:H22	36:5:622:A:N6	2.16	0.44
36:5:989:A:H2'	36:5:990:U:O4'	2.17	0.44
1:6:1218:G:O6	1:6:1444:A:H2'	2.18	0.44
1:6:479:C:H2'	1:6:480:G:O4'	2.18	0.44
1:6:822:U:H2'	1:6:823:G:H5''	2.00	0.44
1:6:841:U:H2'	1:6:842:C:C6	2.53	0.44
1:6:957:G:C5	1:6:958:U:C4	3.05	0.44
37:7:103:A:H2'	37:7:104:A:O4'	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:C0:41:TYR:O	12:C0:45:ALA:N	2.95	0.44
14:C2:52:LEU:O	14:C2:85:LYS:NZ	2.49	0.44
14:C2:66:VAL:HG11	14:C2:71:ILE:HD13	3.11	0.44
16:C4:87:GLY:HA3	16:C4:120:PRO:HG2	2.72	0.44
16:C4:41:ARG:O	16:C4:42:VAL:HG22	2.18	0.44
17:C5:79:HIS:O	17:C5:81:ARG:HG3	2.17	0.44
18:C6:101:SER:HA	18:C6:104:GLU:HB2	1.99	0.44
18:C6:47:LYS:HE3	18:C6:82:ARG:NH1	5.91	0.44
19:C7:18:GLU:OE1	19:C7:19:ARG:NH2	2.48	0.44
20:C8:11:PHE:HA	20:C8:59:GLY:O	2.18	0.44
21:C9:109:GLU:HG2	21:C9:114:VAL:O	3.92	0.44
25:D3:31:LYS:NZ	25:D3:110:LYS:HE3	2.33	0.44
40:L3:306:THR:HG21	40:L3:316:GLU:HG2	4.14	0.44
40:L3:68:HIS:CD2	40:L3:69:LYS:HG3	2.53	0.44
41:L4:141:ARG:N	41:L4:177:ASP:OD1	3.16	0.44
41:L4:203:ARG:HD2	36:5:1383:G:OP1	109.68	0.44
41:L4:44:LYS:HA	41:L4:47:ARG:HD2	2.00	0.44
42:L5:40:HIS:HB3	42:L5:43:LYS:HD2	3.90	0.44
42:L5:75:LEU:O	42:L5:112:LYS:NZ	2.37	0.44
44:L7:168:ILE:O	44:L7:172:ASN:ND2	4.41	0.44
44:L7:89:ILE:HD12	44:L7:214:TRP:CZ3	2.53	0.44
46:L9:86:TYR:CD1	46:L9:151:VAL:HG13	2.53	0.44
46:L9:21:LYS:O	46:L9:22:SER:OG	2.29	0.44
47:M0:100:ASN:OD1	47:M0:118:ALA:HB1	2.17	0.44
47:M0:91:VAL:HG23	47:M0:135:ILE:HA	2.00	0.44
47:M0:210:ILE:HG12	47:M0:217:PHE:CZ	2.53	0.44
36:1:2854:U:OP2	47:M0:3:ARG:NH2	2.51	0.44
51:M5:153:ASP:OD2	51:M5:155:VAL:HG23	4.39	0.44
57:N1:119:ALA:HB3	57:N1:126:VAL:HG13	3.58	0.44
63:N7:47:GLU:N	63:N7:69:LYS:O	2.91	0.44
64:N8:104:THR:HG21	64:N8:112:ILE:HD11	2.05	0.44
67:O1:30:PRO:HG3	67:O1:60:TRP:CZ2	2.52	0.44
36:1:1160:C:N3	68:O2:45:ARG:NH1	2.65	0.44
70:O4:106:LYS:HA	70:O4:106:LYS:HD3	1.79	0.44
71:O5:73:LYS:HD2	71:O5:73:LYS:HA	5.31	0.44
36:1:1844:C:O2	73:O7:9:GLY:HA2	2.18	0.44
4:S2:66:PHE:CD2	4:S2:130:ILE:HG12	4.41	0.44
7:S5:56:ALA:C	7:S5:58:LEU:H	4.26	0.44
8:S6:176:GLN:HG3	8:S6:177:ARG:H	1.83	0.44
10:S8:18:ARG:HG2	10:S8:18:ARG:H	3.86	0.44
10:S8:4:SER:HB3	10:S8:6:ASP:OD2	3.40	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:SR:16:HIS:HB3	34:SR:17:ASN:H	1.71	0.44
34:SR:241:PHE:CG	34:SR:290:VAL:HG12	3.52	0.44
34:SR:84:SER:OG	34:SR:86:ASP:OD2	3.11	0.44
36:1:1139:G:C6	36:1:1140:G:N7	2.86	0.44
36:1:1204:A:H2	36:1:2834:G:N3	2.16	0.44
36:1:1273:A:H2'	36:1:1274:A:H5'	1.99	0.44
36:1:1307:G:H4'	36:1:1308:A:O5'	2.18	0.44
36:1:1447:G:H3'	53:M7:67:ILE:CD1	2.47	0.44
36:1:167:U:H2'	36:1:168:U:C6	2.52	0.44
36:1:1783:U:H2'	36:1:1784:G:H8	1.83	0.44
36:1:1932:A:H5'	36:1:1933:A:OP2	2.17	0.44
36:1:213:A:H2'	36:1:214:G:O4'	2.18	0.44
36:1:2261:G:H21	36:1:2262:A:N6	2.16	0.44
36:1:2401:A:H5'	41:L4:70:ALA:HB2	2.00	0.44
36:1:2729:U:H2'	36:1:2730:G:O4'	2.18	0.44
36:1:3199:G:H5''	50:M4:6:ILE:HG21	2.00	0.44
1:2:1629:G:C5	1:2:1630:U:C4	3.06	0.44
1:2:576:G:H4'	1:2:580:A:C4	2.52	0.44
1:2:77:U:C4	8:S6:170:THR:HG21	2.53	0.44
1:2:987:G:C2	39:L2:249:SER:HB2	2.53	0.44
36:5:1121:U:C4	36:5:1122:U:C4	3.05	0.44
36:5:1741:A:C6	36:5:1742:U:C2	3.06	0.44
55:M9:85:ARG:NH2	36:5:1916:U:O3'	230.54	0.44
36:5:2258:U:H3'	36:5:2259:A:H8	1.82	0.44
36:5:2144:A:H1'	36:5:2281:A:H61	1.83	0.44
36:5:2533:G:N2	36:5:2546:C:H42	2.16	0.44
36:5:2745:G:N2	36:5:2748:A:C8	2.85	0.44
36:5:3106:A:H2'	36:5:3107:U:O4'	2.18	0.44
36:5:664:U:H2'	36:5:665:A:C8	2.53	0.44
36:5:971:G:H2'	36:5:972:A:O4'	2.18	0.44
1:6:1526:A:N1	1:6:1608:U:O2'	2.29	0.44
1:6:778:G:C6	1:6:783:G:C6	3.06	0.44
16:C4:16:VAL:HG23	16:C4:31:THR:HG23	1.99	0.44
17:C5:125:PRO:HG2	17:C5:127:ARG:NH1	4.81	0.44
17:C5:82:ASN:HD22	17:C5:82:ASN:H	1.64	0.44
18:C6:69:VAL:HG21	18:C6:81:ILE:HG12	5.23	0.44
4:S2:156:THR:HB	24:D2:95:PRO:HB3	2.39	0.44
36:1:2554:A:H5''	39:L2:85:GLY:O	2.18	0.44
40:L3:2:SER:O	40:L3:3:HIS:HB3	2.18	0.44
40:L3:68:HIS:NE2	40:L3:69:LYS:HG3	2.33	0.44
40:L3:83:PRO:HB3	40:L3:202:THR:HG23	2.28	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:L4:188:ARG:NE	41:L4:197:ARG:HB3	3.20	0.44
41:L4:215:ILE:HD12	41:L4:218:ALA:HB3	2.00	0.44
41:L4:6:VAL:HG21	41:L4:255:PHE:CZ	2.53	0.44
42:L5:190:ILE:HD11	42:L5:195:LEU:HD22	3.09	0.44
43:L6:131:LYS:HA	43:L6:131:LYS:HD3	4.02	0.44
44:L7:163:LEU:HD21	44:L7:201:PHE:HE1	1.82	0.44
45:L8:163:VAL:HG23	45:L8:166:LEU:HD12	2.00	0.44
45:L8:84:ARG:HH12	45:L8:181:LYS:HZ3	1.65	0.44
51:M5:37:HIS:NE2	51:M5:63:ARG:HB3	2.33	0.44
53:M7:24:VAL:HG13	53:M7:86:LYS:HG2	3.26	0.44
53:M7:48:LEU:O	53:M7:52:LEU:HD22	2.18	0.44
41:L4:281:ILE:HG13	54:M8:125:ASP:CG	2.79	0.44
54:M8:178:ARG:HD3	54:M8:178:ARG:HA	1.48	0.44
36:1:1764:U:H5"	55:M9:43:LYS:HE2	2.00	0.44
56:N0:80:ARG:HB3	56:N0:122:HIS:HB2	1.99	0.44
56:N0:23:LYS:HA	57:N1:146:ASN:OD1	3.56	0.44
62:N6:43:TYR:CD1	62:N6:126:LEU:HA	2.53	0.44
66:O0:15:ALA:O	66:O0:19:LYS:HG2	2.18	0.44
67:O1:52:ALA:HA	67:O1:53:PRO:HD3	1.88	0.44
71:O5:66:VAL:HG12	71:O5:80:LEU:HD11	2.57	0.44
2:S0:67:ILE:HA	2:S0:68:PRO:HD3	1.95	0.44
3:S1:134:VAL:O	3:S1:218:LEU:HD22	2.18	0.44
3:S1:175:GLU:HG3	3:S1:187:LYS:HD3	5.37	0.44
3:S1:190:PRO:HG2	3:S1:192:VAL:HG23	2.19	0.44
4:S2:101:VAL:HG22	4:S2:115:ILE:HG12	1.99	0.44
5:S3:23:GLU:OE2	12:C0:61:TRP:NE1	2.76	0.44
7:S5:35:GLN:HG2	7:S5:39:GLU:OE1	2.18	0.44
10:S8:8:ARG:NH2	10:S8:19:ALA:O	2.49	0.44
11:S9:115:LYS:HD2	11:S9:115:LYS:HA	1.87	0.44
11:S9:36:LEU:HA	11:S9:126:ARG:NH2	2.33	0.44
11:S9:149:ARG:H	11:S9:149:ARG:HG2	1.53	0.44
35:SM:33:LYS:HB2	35:SM:33:LYS:HE3	4.70	0.44
35:SM:84:LYS:O	35:SM:85:SER:HB3	2.18	0.44
34:SR:282:SER:O	34:SR:286:GLU:HG3	2.18	0.44
36:1:1720:U:OP2	55:M9:110:ARG:NH1	2.51	0.44
36:1:2947:G:H4'	36:1:2947:G:OP2	2.17	0.44
36:1:3331:U:H2'	36:1:3332:U:O4'	2.18	0.44
36:1:3386:G:H2'	36:1:3387:U:C6	2.53	0.44
1:2:1029:U:O2'	1:2:1030:A:H5'	2.17	0.44
1:2:1524:A:C2	1:2:1590:G:H1'	2.53	0.44
1:2:66:U:C4	8:S6:173:PRO:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:1638:A:N1	36:5:1736:G:O2'	2.47	0.44
36:5:2396:G:N2	36:5:2985:C:C2	2.86	0.44
36:5:343:U:O2	36:5:1439:U:H1'	2.17	0.44
36:5:568:G:O5'	36:5:568:G:H8	2.01	0.44
1:6:1240:U:H3	1:6:1242:A:H5''	1.83	0.44
1:6:40:A:H2'	1:6:41:A:O4'	2.18	0.44
1:6:504:U:O4	1:6:505:A:N6	2.50	0.44
1:6:832:U:H2'	1:6:833:U:O4'	2.18	0.44
38:8:76:C:H2'	38:8:77:A:O4'	2.17	0.44
13:C1:118:GLN:HG3	13:C1:121:ASP:OD2	2.18	0.44
17:C5:60:LEU:HA	17:C5:76:VAL:HG21	1.99	0.44
20:C8:38:VAL:HG11	20:C8:73:MET:SD	2.57	0.44
22:D0:23:ARG:O	22:D0:117:VAL:HG12	2.18	0.44
26:D4:111:LYS:HD3	26:D4:114:ARG:NH2	5.01	0.44
11:S9:29:LYS:HA	32:E0:40:TYR:CE2	2.58	0.44
39:L2:97:ASN:HB2	39:L2:100:ASN:HD22	2.94	0.44
39:L2:169:ILE:O	39:L2:171:GLY:N	3.97	0.44
41:L4:192:GLY:HA2	41:L4:195:ARG:HG3	3.64	0.44
41:L4:60:THR:HG23	36:5:364:G:OP1	129.70	0.44
42:L5:78:ALA:HB3	42:L5:105:ILE:HG12	2.00	0.44
44:L7:162:PRO:O	44:L7:164:SER:N	2.88	0.44
46:L9:92:TYR:N	46:L9:92:TYR:CD1	2.83	0.44
48:M1:166:LYS:O	48:M1:168:ASP:N	3.17	0.44
49:M3:172:LEU:O	49:M3:176:GLU:HB3	3.28	0.44
49:M3:185:LYS:HG3	49:M3:188:ARG:NH1	4.97	0.44
49:M3:89:TYR:O	49:M3:92:THR:OG1	2.93	0.44
50:M4:112:LEU:HD22	50:M4:116:GLU:HG2	3.73	0.44
51:M5:104:GLU:HG2	51:M5:160:GLU:HG2	2.37	0.44
57:N1:9:SER:O	57:N1:55:LYS:NZ	2.93	0.44
58:N2:59:ASP:HB3	58:N2:62:VAL:HG12	2.00	0.44
59:N3:28:ASN:OD1	59:N3:111:GLY:HA3	3.62	0.44
59:N3:32:ARG:O	59:N3:32:ARG:NH1	7.66	0.44
62:N6:40:ARG:O	62:N6:44:GLY:N	2.51	0.44
63:N7:83:THR:HG23	70:O4:93:PHE:CE1	5.74	0.44
63:N7:88:ASP:HB3	63:N7:121:ARG:NH2	2.31	0.44
74:O8:11:PHE:CD1	74:O8:12:LEU:HD22	2.62	0.44
74:O8:58:ASP:HB3	74:O8:61:LYS:HG2	4.24	0.44
2:S0:125:ASP:O	2:S0:129:ASP:HB2	3.95	0.44
3:S1:194:ASN:N	3:S1:194:ASN:OD1	2.49	0.44
5:S3:149:ALA:O	5:S3:150:MET:HG2	2.18	0.44
6:S4:221:ARG:O	6:S4:224:ASN:N	3.12	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:S4:42:LEU:HD12	6:S4:109:PHE:HD2	5.39	0.44
7:S5:94:THR:HG22	7:S5:114:ILE:HG21	1.98	0.44
7:S5:143:ARG:HD2	30:D8:57:MET:SD	2.58	0.44
7:S5:56:ALA:O	7:S5:58:LEU:N	3.51	0.44
8:S6:142:ARG:O	8:S6:145:PHE:N	3.52	0.44
9:S7:39:ARG:HG3	9:S7:40:PRO:HD3	1.99	0.44
35:SM:41:SER:O	35:SM:43:ASP:N	2.43	0.44
34:SR:145:LEU:H	34:SR:145:LEU:HD23	3.99	0.44
34:SR:32:LEU:HD11	34:SR:44:SER:HB2	2.45	0.44
36:1:1307:G:C2	36:1:1308:A:C2	3.06	0.43
36:1:1699:A:H2'	36:1:1700:G:H8	1.83	0.43
36:1:692:A:C4	36:1:693:A:C8	3.06	0.43
36:1:802:C:H2'	36:1:803:C:H6	1.84	0.43
1:2:1588:G:H2'	1:2:1589:C:H5'	2.00	0.43
1:2:387:A:H3'	1:2:402:C:H5	1.83	0.43
1:2:406:U:O3'	8:S6:94:ARG:NH2	2.35	0.43
1:2:411:C:H2'	1:2:412:A:O4'	2.17	0.43
1:2:700:C:H42	1:2:738:G:H22	1.65	0.43
1:2:700:C:N4	1:2:738:G:H22	2.16	0.43
36:5:1081:U:H1'	36:5:1082:U:OP2	2.18	0.43
36:5:1348:U:O4'	36:5:1355:A:N6	2.51	0.43
36:5:2566:C:H42	36:5:2575:G:H1	1.67	0.43
36:5:532:A:O2'	36:5:533:A:H5'	2.18	0.43
36:5:686:G:C6	36:5:687:U:C2	3.06	0.43
54:M8:90:ASP:HA	36:5:785:G:O6	139.17	0.43
1:6:1274:C:O2	1:6:1274:C:H2'	2.18	0.43
1:6:503:G:H2'	1:6:504:U:H6	1.83	0.43
1:6:988:A:C2	1:6:989:U:H1'	2.53	0.43
12:C0:58:GLN:HB3	12:C0:65:TYR:HB2	2.82	0.43
12:C0:71:GLU:H	12:C0:71:GLU:HG2	3.72	0.43
14:C2:125:ASN:O	14:C2:126:TRP:HD1	2.01	0.43
17:C5:32:ASP:N	17:C5:32:ASP:OD2	2.51	0.43
19:C7:41:ILE:HD13	19:C7:47:ARG:HA	2.00	0.43
21:C9:108:LEU:HD22	21:C9:113:ILE:HG13	2.44	0.43
21:C9:34:VAL:HG23	21:C9:53:TRP:NE1	2.33	0.43
22:D0:65:ILE:HG21	31:D9:43:PHE:CE1	2.53	0.43
23:D1:36:VAL:HG11	23:D1:78:LEU:HD13	1.99	0.43
25:D3:76:LEU:HB2	25:D3:79:ASN:ND2	2.33	0.43
1:2:525:A:H4'	26:D4:89:TYR:CG	2.53	0.43
30:D8:26:THR:HG22	30:D8:28:VAL:HG23	1.99	0.43
33:E1:144:CYS:CB	33:E1:147:VAL:HG22	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:L3:194:TRP:CE2	40:L3:198:HIS:CE1	3.06	0.43
41:L4:141:ARG:NH1	41:L4:180:LYS:HE2	4.47	0.43
41:L4:233:LEU:HD23	41:L4:233:LEU:HA	1.93	0.43
42:L5:40:HIS:HD2	42:L5:42:ALA:N	2.09	0.43
43:L6:19:LYS:O	43:L6:21:THR:N	2.50	0.43
45:L8:252:ASN:C	45:L8:254:ASP:H	2.21	0.43
46:L9:23:ARG:HH21	46:L9:42:ASP:HA	1.83	0.43
50:M4:22:LEU:HD12	50:M4:32:LEU:HG	4.36	0.43
51:M5:150:TRP:O	51:M5:153:ASP:HB2	2.44	0.43
51:M5:154:PRO:O	51:M5:157:LYS:HB2	2.18	0.43
52:M6:148:LYS:HB2	52:M6:149:TYR:H	1.54	0.43
53:M7:88:VAL:O	53:M7:92:GLN:HG2	2.18	0.43
62:N6:58:VAL:HB	62:N6:63:LYS:O	2.18	0.43
67:O1:44:MET:O	67:O1:46:THR:N	3.34	0.43
71:O5:10:ARG:NH2	38:8:65:A:O3'	32.76	0.43
36:1:2554:A:H62	79:Q3:62:LYS:HZ2	1.65	0.43
39:L2:112:ILE:HD11	79:Q3:79:VAL:HG11	5.05	0.43
79:Q3:8:VAL:HB	79:Q3:11:THR:HG22	1.99	0.43
2:S0:184:LEU:HD13	23:D1:43:GLY:O	2.18	0.43
2:S0:4:PRO:HB2	2:S0:5:ALA:H	1.63	0.43
4:S2:139:ILE:CD1	4:S2:191:ALA:HB1	2.60	0.43
4:S2:239:PRO:HA	4:S2:242:ILE:HB	2.73	0.43
10:S8:59:ARG:O	10:S8:60:ILE:HG13	2.50	0.43
1:2:258:C:H5'	10:S8:75:LYS:HB2	1.99	0.43
34:SR:21:THR:O	34:SR:291:SER:HB3	2.18	0.43
36:1:1016:C:H2'	36:1:1017:C:H5'	2.00	0.43
36:1:1108:U:H2'	36:1:1109:U:C6	2.52	0.43
36:1:1306:G:C6	52:M6:62:THR:HA	2.53	0.43
36:1:1632:A:H2'	36:1:1633:C:C6	2.53	0.43
36:1:1658:G:C5	36:1:1659:U:C5	3.07	0.43
36:1:1495:U:H5	36:1:1835:A:N1	2.15	0.43
36:1:2860:U:C6	36:1:2938:G:H4'	2.52	0.43
36:1:3174:A:C5	36:1:3279:A:H1'	2.53	0.43
36:1:508:U:H2'	36:1:509:U:H6	1.82	0.43
36:1:587:U:H2'	36:1:588:G:H5'	1.99	0.43
36:1:901:G:H2'	36:1:902:G:H8	1.83	0.43
36:1:997:A:H2'	36:1:998:A:O4'	2.18	0.43
1:2:1344:A:H4'	1:2:1345:A:OP1	2.17	0.43
1:2:1559:A:H4'	1:2:1559:A:OP1	2.18	0.43
1:2:1628:U:H2'	1:2:1629:G:C8	2.53	0.43
1:2:46:A:O2'	1:2:432:G:N2	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:686:C:H2'	1:2:687:G:C8	2.53	0.43
36:5:1301:A:H4'	36:5:1302:A:O5'	2.18	0.43
36:5:1615:C:H2'	36:5:1616:U:C6	2.53	0.43
36:5:884:A:C8	36:5:2139:A:C8	3.07	0.43
36:5:2992:U:H2'	36:5:2993:G:O4'	2.18	0.43
36:5:3131:U:H2'	36:5:3132:C:C6	2.52	0.43
36:5:3194:C:C2	36:5:3197:G:N2	2.84	0.43
36:5:3233:C:H2'	36:5:3234:A:C8	2.53	0.43
36:5:80:G:H2'	36:5:81:C:C6	2.53	0.43
1:6:1636:C:C4	1:6:1765:A:N1	2.87	0.43
1:6:197:A:H2'	1:6:198:A:H8	1.83	0.43
1:6:461:G:H2'	1:6:462:G:C8	2.53	0.43
1:6:69:G:H2'	1:6:70:C:O4'	2.18	0.43
6:S4:187:ARG:NH1	1:6:753:A:N7	373.29	0.43
24:D2:57:ARG:HG2	1:6:863:A:O5'	351.09	0.43
18:C6:127:LYS:NZ	1:6:1604:U:OP2	384.28	0.43
19:C7:21:TYR:N	19:C7:22:PRO:HD2	2.32	0.43
19:C7:24:LEU:HD23	19:C7:34:LEU:HD13	4.91	0.43
20:C8:33:THR:C	20:C8:35:ILE:H	2.21	0.43
26:D4:53:ASP:OD2	26:D4:96:LEU:HD21	2.18	0.43
39:L2:180:LEU:HD23	39:L2:180:LEU:HA	1.71	0.43
42:L5:179:ARG:HA	42:L5:179:ARG:HD3	2.45	0.43
43:L6:31:ARG:NH2	43:L6:81:ALA:O	2.51	0.43
47:M0:119:TRP:CG	47:M0:120:GLY:N	2.86	0.43
48:M1:82:ARG:HD2	48:M1:112:LEU:HB2	4.32	0.43
49:M3:159:VAL:HG11	64:N8:142:GLY:O	2.17	0.43
52:M6:98:ALA:HA	52:M6:101:ARG:NH1	2.96	0.43
57:N1:69:LYS:HB3	36:5:2737:C:OP1	229.17	0.43
57:N1:83:ARG:HG2	57:N1:84:TYR:N	2.32	0.43
64:N8:64:GLN:HB2	64:N8:67:HIS:CD2	2.52	0.43
70:O4:107:GLU:HG2	70:O4:110:GLU:OE2	6.71	0.43
70:O4:8:ARG:HD3	36:5:1606:U:C4	138.76	0.43
73:O7:58:THR:O	73:O7:61:THR:HG23	2.18	0.43
74:O8:58:ASP:OD2	74:O8:61:LYS:N	2.29	0.43
4:S2:83:ILE:HG12	4:S2:100:ALA:HA	2.00	0.43
4:S2:225:LEU:HA	4:S2:225:LEU:HD23	1.76	0.43
4:S2:230:TRP:NE1	24:D2:68:ARG:HB2	4.53	0.43
5:S3:119:ALA:O	5:S3:123:VAL:HG23	2.24	0.43
6:S4:246:LEU:HB2	6:S4:251:GLU:HG3	2.00	0.43
6:S4:73:ASP:OD1	6:S4:89:VAL:N	2.35	0.43
7:S5:112:ARG:HH11	18:C6:43:ILE:HD11	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:S5:146:THR:HG23	7:S5:221:ALA:HA	1.99	0.43
8:S6:145:PHE:HB2	8:S6:147:LEU:HG	1.98	0.43
8:S6:85:ARG:HB3	8:S6:87:ARG:HH21	4.14	0.43
10:S8:21:PHE:HD1	10:S8:22:ARG:HG2	3.95	0.43
11:S9:96:VAL:HG23	11:S9:99:LEU:HD22	6.21	0.43
34:SR:135:THR:OG1	34:SR:136:ILE:N	2.49	0.43
36:1:1103:A:H1'	36:1:1104:G:P	2.59	0.43
36:1:1369:A:OP1	64:N8:21:ARG:NH1	2.52	0.43
36:1:1658:G:C6	36:1:1659:U:C4	3.07	0.43
36:1:2413:A:H2'	36:1:2414:G:H8	1.83	0.43
36:1:505:G:C6	36:1:506:U:C4	3.07	0.43
36:1:686:G:P	49:M3:39:ARG:HH21	2.41	0.43
1:2:1076:A:O3'	28:D6:13:LYS:NZ	2.42	0.43
1:2:357:G:H2'	1:2:358:U:O4'	2.17	0.43
1:2:512:A:H2'	1:2:513:U:C6	2.53	0.43
1:2:642:G:H2'	1:2:643:G:C8	2.53	0.43
37:3:22:A:H2'	37:3:23:A:C8	2.53	0.43
37:3:87:G:C2	37:3:88:G:C8	3.06	0.43
38:4:36:G:N2	38:4:37:A:N1	2.66	0.43
36:5:953:G:O2'	36:5:1116:G:H5'	2.18	0.43
36:5:1275:C:H2'	36:5:1276:U:O4'	2.18	0.43
70:O4:4:ARG:HD2	36:5:1485:G:N2	151.66	0.43
36:5:1812:G:H4'	36:5:1812:G:OP1	2.18	0.43
36:5:2287:C:C2	36:5:2298:U:O4'	2.71	0.43
36:5:2665:U:H4'	36:5:2666:C:OP1	2.18	0.43
36:5:278:U:H2'	36:5:279:U:C6	2.53	0.43
36:5:293:C:H2'	36:5:294:U:O4'	2.18	0.43
36:5:2948:C:H2'	36:5:2949:U:C6	2.53	0.43
53:M7:69:ARG:NH2	36:5:2992:U:H1'	191.91	0.43
36:5:523:A:N6	36:5:570:A:C2	2.86	0.43
36:5:656:A:H2'	36:5:657:A:C8	2.53	0.43
36:5:80:G:H2'	36:5:81:C:H6	1.84	0.43
36:5:945:C:H1'	36:5:1407:A:H1'	2.00	0.43
1:6:1029:U:O2'	1:6:1030:A:H5'	2.18	0.43
1:6:1561:U:H2'	1:6:1562:G:H8	1.84	0.43
1:6:1578:U:H2'	1:6:1579:U:C6	2.53	0.43
18:C6:122:ARG:HB3	1:6:1584:G:C8	395.49	0.43
1:6:1620:C:O2'	1:6:1621:U:OP1	2.33	0.43
1:6:587:C:H2'	1:6:588:U:O4'	2.18	0.43
16:C4:37:GLU:HA	1:6:895:G:O2'	259.18	0.43
12:C0:21:VAL:HG12	12:C0:22:VAL:H	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:C1:5:LEU:HD23	13:C1:7:VAL:N	6.79	0.43
17:C5:52:LYS:N	17:C5:53:PRO:HD2	3.44	0.43
18:C6:66:ARG:NH2	18:C6:68:ARG:HD2	2.33	0.43
1:2:1559:A:C6	20:C8:134:ARG:HD2	2.53	0.43
20:C8:36:LYS:NZ	1:6:1568:C:OP1	334.16	0.43
27:D5:102:THR:HB	27:D5:103:ARG:H	1.59	0.43
27:D5:87:GLY:O	27:D5:89:ILE:N	2.42	0.43
31:D9:31:ILE:HA	31:D9:31:ILE:HD13	2.20	0.43
32:E0:3:LYS:HB3	32:E0:3:LYS:HE3	1.89	0.43
33:E1:148:TYR:HD1	33:E1:148:TYR:HA	1.99	0.43
40:L3:13:HIS:HB3	40:L3:16:PHE:HD1	1.97	0.43
41:L4:181:VAL:HG11	41:L4:224:GLY:CA	3.38	0.43
43:L6:5:LYS:HA	43:L6:5:LYS:HE3	2.00	0.43
43:L6:98:VAL:HA	43:L6:101:PHE:HD2	1.82	0.43
44:L7:179:LEU:H	44:L7:179:LEU:HD22	1.83	0.43
45:L8:160:ILE:HG12	45:L8:160:ILE:H	1.50	0.43
45:L8:160:ILE:HG22	45:L8:164:VAL:HG13	2.00	0.43
47:M0:16:PRO:HG3	47:M0:128:ARG:NH1	2.33	0.43
48:M1:109:HIS:HE1	48:M1:122:ILE:HA	1.82	0.43
49:M3:127:PRO:HG2	49:M3:131:LYS:HD2	2.00	0.43
52:M6:110:PRO:O	52:M6:111:PRO:C	2.98	0.43
52:M6:16:VAL:HG23	52:M6:42:ASN:O	2.26	0.43
57:N1:120:LYS:O	57:N1:122:GLN:N	2.51	0.43
57:N1:17:ARG:HD2	57:N1:17:ARG:HA	1.98	0.43
57:N1:73:GLY:HA2	57:N1:89:LEU:O	2.32	0.43
59:N3:75:PRO:HB2	59:N3:103:ALA:O	2.18	0.43
61:N5:45:LYS:HB3	71:O5:75:TYR:HD2	3.51	0.43
64:N8:93:SER:OG	64:N8:93:SER:O	2.36	0.43
68:O2:4:LEU:H	68:O2:4:LEU:HD22	4.65	0.43
76:Q0:110:CYS:SG	76:Q0:111:ARG:N	2.91	0.43
2:S0:10:THR:OG1	2:S0:13:ASP:OD2	2.36	0.43
2:S0:58:VAL:O	2:S0:62:ARG:HG2	5.22	0.43
3:S1:209:ASN:HB3	3:S1:211:HIS:NE2	2.34	0.43
4:S2:157:LYS:HD3	4:S2:170:ILE:HG23	2.00	0.43
6:S4:86:PHE:CE1	6:S4:87:MET:HG2	2.54	0.43
10:S8:21:PHE:CE1	10:S8:22:ARG:HD3	2.53	0.43
10:S8:3:ILE:HB	10:S8:30:GLY:O	2.28	0.43
36:1:1329:U:H1'	36:1:1330:A:OP1	2.18	0.43
36:1:1483:G:C8	36:1:1485:G:C8	3.06	0.43
36:1:1823:A:H2'	36:1:1824:U:C6	2.54	0.43
36:1:372:A:H2'	36:1:373:A:O4'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:806:A:C4	36:1:936:A:C2	3.06	0.43
36:1:952:A:N3	36:1:1114:U:O2'	2.37	0.43
1:2:1002:G:H2'	1:2:1003:A:H5'	2.00	0.43
1:2:1105:C:H41	25:D3:4:GLY:HA3	1.84	0.43
1:2:1498:G:O2'	1:2:1499:G:H5'	2.18	0.43
1:2:1514:U:H5''	1:2:1515:A:O4'	2.19	0.43
1:2:566:C:H2'	1:2:567:A:O4'	2.19	0.43
1:2:763:G:C6	1:2:764:U:C4	3.06	0.43
1:2:927:C:O2'	16:C4:124:ASP:O	2.36	0.43
36:5:1560:G:O2'	36:5:1561:G:OP1	2.31	0.43
41:L4:217:LYS:NZ	36:5:210:U:O2	66.19	0.43
36:5:1125:U:O2'	36:5:2643:A:N1	2.49	0.43
54:M8:43:PRO:HB2	36:5:728:G:H5''	190.25	0.43
36:5:913:A:H5''	36:5:914:A:N7	2.34	0.43
1:6:1405:G:H2'	1:6:1406:A:C8	2.54	0.43
35:SM:71:ASN:HB3	1:6:1460:A:O4'	329.94	0.43
20:C8:134:ARG:NH1	1:6:1546:G:OP2	358.77	0.43
1:6:567:A:C2	1:6:583:C:H1'	2.53	0.43
1:6:955:A:H2'	1:6:956:C:O4'	2.18	0.43
14:C2:83:GLU:C	14:C2:85:LYS:H	4.86	0.43
16:C4:91:THR:O	16:C4:92:LYS:HG2	2.19	0.43
28:D6:4:LYS:NZ	28:D6:5:ARG:HH22	5.13	0.43
30:D8:16:LEU:HD12	30:D8:27:GLN:OE1	7.17	0.43
39:L2:241:ARG:HG2	39:L2:241:ARG:H	2.26	0.43
40:L3:290:ASP:OD2	40:L3:292:ALA:N	4.91	0.43
40:L3:37:ARG:O	40:L3:186:GLY:HA2	2.18	0.43
42:L5:279:LYS:HD3	42:L5:279:LYS:HA	2.22	0.43
44:L7:33:ARG:NH2	36:5:595:G:OP1	232.65	0.43
49:M3:6:ASN:O	49:M3:7:LEU:HD23	2.82	0.43
49:M3:93:ILE:O	49:M3:119:TYR:OH	2.21	0.43
52:M6:119:VAL:O	52:M6:121:PRO:HD3	2.17	0.43
53:M7:92:GLN:HA	53:M7:95:LEU:HB2	2.00	0.43
54:M8:96:PHE:CE1	54:M8:114:ILE:HA	2.78	0.43
55:M9:8:LYS:HG3	55:M9:22:VAL:CG2	2.48	0.43
56:N0:23:LYS:HE3	56:N0:25:PHE:HZ	1.82	0.43
59:N3:10:LYS:HZ3	59:N3:56:ASP:CG	2.22	0.43
8:S6:70:PRO:HG3	60:N4:2:LYS:HE3	4.71	0.43
60:N4:59:HIS:HB3	60:N4:61:LYS:HE2	2.01	0.43
62:N6:12:ARG:O	62:N6:12:ARG:HD2	5.02	0.43
62:N6:55:GLU:HB3	62:N6:108:LYS:H	1.84	0.43
63:N7:81:LEU:HA	63:N7:81:LEU:HD22	2.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:97:SER:H	63:N7:100:THR:HB	1.83	0.43
51:M5:9:GLU:HG3	72:O6:44:VAL:HG11	2.87	0.43
72:O6:54:GLU:O	72:O6:58:ILE:HG23	2.30	0.43
36:1:2138:A:C5	73:O7:3:LYS:HB3	2.54	0.43
78:Q2:38:GLN:NE2	78:Q2:38:GLN:O	2.46	0.43
2:S0:179:ARG:HD3	2:S0:183:ARG:CZ	3.42	0.43
4:S2:88:LYS:HB3	4:S2:95:ARG:HB3	2.97	0.43
4:S2:98:PHE:CZ	35:SM:116:GLU:HG3	2.54	0.43
5:S3:102:ALA:HB1	5:S3:173:ARG:CG	2.94	0.43
5:S3:27:ARG:HB3	12:C0:58:GLN:NE2	2.71	0.43
7:S5:177:ILE:HA	7:S5:180:ARG:NH1	2.34	0.43
9:S7:138:LYS:HB2	24:D2:54:ASP:HB3	3.55	0.43
9:S7:153:LEU:HD22	9:S7:184:GLU:HB2	2.00	0.43
9:S7:45:SER:O	9:S7:61:PHE:N	3.00	0.43
11:S9:13:SER:HB2	11:S9:47:PHE:CD1	2.53	0.43
34:SR:175:ASP:N	34:SR:175:ASP:OD1	2.56	0.43
34:SR:22:SER:OG	34:SR:70:ASP:HA	3.26	0.43
36:1:1724:U:H4'	36:1:1725:C:OP1	2.19	0.43
36:1:1920:U:N3	36:1:1930:A:C2	2.87	0.43
36:1:207:U:H2'	36:1:208:C:C6	2.54	0.43
36:1:2113:A:N7	36:1:2114:C:C4	2.86	0.43
36:1:3028:G:H2'	36:1:3029:A:O4'	2.19	0.43
36:1:3163:A:N1	36:1:3164:C:N4	2.66	0.43
36:1:620:U:C4	36:1:622:A:C6	3.06	0.43
1:2:1017:U:H2'	1:2:1018:U:C6	2.53	0.43
1:2:330:G:OP2	10:S8:172:ARG:NH1	2.48	0.43
1:2:712:G:H1	1:2:713:A:H62	1.66	0.43
1:2:729:G:C6	1:2:730:G:H1'	2.53	0.43
36:5:1257:C:H2'	36:5:1258:U:O4'	2.18	0.43
36:5:1481:A:H2'	36:5:1481:A:N3	2.34	0.43
36:5:1765:U:OP1	36:5:1765:U:H4'	2.18	0.43
36:5:21:G:OP2	38:8:36:G:N1	2.51	0.43
36:5:2426:U:H2'	36:5:2427:U:C6	2.54	0.43
36:5:2775:U:H2'	36:5:2776:C:H6	1.82	0.43
36:5:853:G:H2'	36:5:854:G:O4'	2.18	0.43
1:6:1292:G:C6	1:6:1293:U:C4	3.07	0.43
1:6:21:U:H2'	1:6:22:A:C8	2.53	0.43
1:6:224:C:H2'	1:6:225:A:C8	2.43	0.43
1:6:791:A:H2'	1:6:792:U:O4'	2.18	0.43
14:C2:91:VAL:HG22	14:C2:92:ALA:H	4.39	0.43
15:C3:27:LYS:HD2	15:C3:28:LEU:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:C5:33:PHE:O	17:C5:36:LEU:HD23	2.19	0.43
18:C6:31:VAL:HG11	18:C6:81:ILE:HD13	2.89	0.43
20:C8:127:HIS:NE2	20:C8:133:VAL:HG11	2.33	0.43
20:C8:89:GLN:C	20:C8:91:ASP:H	4.60	0.43
1:2:1199:G:O6	22:D0:67:THR:HG23	2.19	0.43
22:D0:58:LEU:HD12	22:D0:88:LYS:HD2	2.00	0.43
23:D1:71:ARG:HB2	23:D1:83:TRP:CD1	5.27	0.43
25:D3:23:ARG:HB3	25:D3:29:TYR:CE1	2.59	0.43
32:E0:41:THR:HA	32:E0:45:VAL:HB	2.01	0.43
40:L3:252:ILE:HG22	36:5:2394:G:H5'	216.67	0.43
41:L4:22:LEU:HA	41:L4:23:PRO:HD3	1.81	0.43
42:L5:65:ILE:HA	42:L5:74:VAL:HA	2.88	0.43
44:L7:173:LEU:HD12	44:L7:173:LEU:HA	2.05	0.43
45:L8:109:LEU:HA	45:L8:109:LEU:HD13	2.27	0.43
48:M1:137:ARG:HB2	48:M1:137:ARG:HE	1.39	0.43
48:M1:133:ARG:HH12	48:M1:154:THR:HA	1.83	0.43
49:M3:168:ARG:NH1	49:M3:172:LEU:HD11	2.34	0.43
49:M3:58:VAL:N	49:M3:70:ARG:O	2.88	0.43
50:M4:113:THR:HG22	50:M4:115:PHE:H	1.83	0.43
51:M5:99:ARG:HH21	51:M5:166:ALA:HB3	3.34	0.43
52:M6:120:VAL:HG12	52:M6:122:GLN:HG3	1.99	0.43
52:M6:187:GLU:HB3	52:M6:192:LYS:NZ	5.54	0.43
54:M8:60:PRO:HG2	54:M8:142:GLY:CA	3.63	0.43
54:M8:86:THR:HA	54:M8:105:ARG:O	2.19	0.43
59:N3:13:ILE:HG12	59:N3:53:SER:HB2	5.61	0.43
59:N3:94:TYR:HE2	60:N4:19:THR:OG1	2.01	0.43
62:N6:70:ILE:HG12	62:N6:80:VAL:HG21	2.00	0.43
63:N7:119:GLU:O	63:N7:123:GLN:HG3	2.96	0.43
63:N7:17:ARG:HA	70:O4:74:ARG:HA	2.00	0.43
70:O4:21:LYS:HG2	70:O4:22:VAL:N	3.72	0.43
72:O6:33:ALA:O	72:O6:34:SER:OG	2.25	0.43
73:O7:19:CYS:O	73:O7:23:GLY:N	2.40	0.43
2:S0:57:LEU:HD11	2:S0:176:LEU:HB2	2.01	0.43
5:S3:160:SER:O	1:6:1420:C:O2'	414.14	0.43
5:S3:162:GLN:N	5:S3:163:PRO:HD2	2.57	0.43
5:S3:71:LEU:HD22	5:S3:75:LYS:HE2	2.00	0.43
5:S3:93:ASP:N	5:S3:93:ASP:OD2	2.51	0.43
6:S4:26:CYS:SG	11:S9:2:PRO:HB2	2.59	0.43
8:S6:176:GLN:HG3	8:S6:177:ARG:N	2.51	0.43
1:2:260:U:O4	10:S8:42:ARG:HA	2.18	0.43
11:S9:99:LEU:CD1	11:S9:100:LYS:H	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:1235:U:H4'	36:1:1236:G:H5'	1.98	0.43
36:1:1472:U:H2'	36:1:1473:G:H8	1.83	0.43
36:1:2407:C:H2'	36:1:2408:U:C6	2.54	0.43
36:1:2946:A:C5'	36:1:2947:G:H5'	2.49	0.43
36:1:3298:C:H2'	36:1:3299:A:O4'	2.19	0.43
36:1:595:G:H2'	36:1:596:C:H6	1.82	0.43
36:1:603:A:H2'	36:1:604:G:O4'	2.18	0.43
36:1:901:G:C4	36:1:902:G:C8	3.07	0.43
1:2:1355:C:O2	1:2:1368:G:N2	2.42	0.43
1:2:1352:G:H1	1:2:1373:C:N4	2.16	0.43
1:2:1746:A:H2'	1:2:1747:G:O4'	2.18	0.43
1:2:1762:A:H1'	1:2:1783:C:H5'	2.00	0.43
1:2:649:U:O2'	1:2:650:U:O5'	2.36	0.43
36:5:1047:A:C6	36:5:1048:A:C6	3.07	0.43
36:5:1141:C:H2'	36:5:1142:G:O4'	2.18	0.43
36:5:1479:U:H2'	36:5:1480:G:H5'	2.01	0.43
36:5:1658:G:C4	36:5:1796:G:C6	3.06	0.43
39:L2:227:ARG:NH2	36:5:2155:G:O2'	205.30	0.43
36:5:215:G:N2	36:5:225:C:O2	2.48	0.43
49:M3:45:LYS:HG3	36:5:241:G:H4'	44.05	0.43
36:5:3153:U:H1'	36:5:3154:C:C2	2.54	0.43
36:5:612:U:H2'	36:5:613:G:H8	1.83	0.43
1:6:1058:U:H1'	1:6:1059:U:H5''	2.00	0.43
29:D7:17:ARG:HD2	1:6:1070:C:H4'	369.67	0.43
24:D2:71:LYS:HD2	1:6:1098:U:O2'	379.45	0.43
1:6:546:U:H2'	1:6:547:U:H6	1.84	0.43
16:C4:54:GLU:CD	1:6:901:G:H22	282.68	0.43
38:8:68:G:C6	38:8:69:U:C4	3.06	0.43
14:C2:67:THR:C	14:C2:69:ALA:H	2.35	0.43
15:C3:20:ARG:HD3	24:D2:56:HIS:NE2	4.76	0.43
15:C3:52:VAL:HG22	15:C3:55:ARG:NH2	2.33	0.43
15:C3:61:THR:O	15:C3:62:GLN:HG3	2.39	0.43
7:S5:69:PHE:CD2	18:C6:50:GLU:HG3	2.54	0.43
20:C8:136:GLN:HG2	20:C8:136:GLN:H	1.38	0.43
21:C9:6:VAL:HB	21:C9:14:PHE:CE1	3.03	0.43
22:D0:106:ILE:HA	22:D0:106:ILE:HD12	1.86	0.43
24:D2:103:ILE:HD13	24:D2:126:LEU:HB2	2.00	0.43
24:D2:14:ILE:HD11	24:D2:38:LEU:HD21	2.00	0.43
1:2:150:U:P	26:D4:123:LYS:HZ3	2.40	0.43
26:D4:27:VAL:HG12	26:D4:29:HIS:HD2	3.06	0.43
1:2:526:A:P	26:D4:93:ARG:HH21	2.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:D6:12:LYS:HB3	28:D6:13:LYS:H	4.11	0.43
32:E0:10:ARG:O	32:E0:12:GLY:N	3.27	0.43
41:L4:26:PHE:CD1	41:L4:130:ALA:HB2	2.53	0.43
42:L5:104:LEU:HD11	42:L5:108:ARG:HH21	1.83	0.43
42:L5:34:LYS:HE3	57:N1:30:TYR:CE1	3.95	0.43
44:L7:90:LYS:HB3	44:L7:133:TYR:HB3	2.21	0.43
47:M0:43:VAL:HA	47:M0:139:ARG:NH2	2.32	0.43
48:M1:25:GLU:OE1	48:M1:29:ARG:HB3	2.85	0.43
48:M1:6:GLN:HB3	48:M1:7:ASN:H	2.02	0.43
51:M5:73:ARG:O	51:M5:75:VAL:HG22	2.68	0.43
51:M5:74:PRO:HB2	51:M5:75:VAL:H	1.61	0.43
53:M7:60:PHE:HB3	53:M7:64:ASN:HB3	2.00	0.43
55:M9:168:ALA:HB1	55:M9:172:ARG:HD2	1.99	0.43
61:N5:57:LEU:HD21	61:N5:90:ALA:HB2	2.01	0.43
62:N6:110:HIS:O	62:N6:110:HIS:ND1	2.51	0.43
62:N6:124:GLY:O	62:N6:126:LEU:N	4.29	0.43
65:N9:5:LYS:HD3	65:N9:8:THR:HG22	2.01	0.43
36:1:1456:A:N6	67:O1:64:VAL:HG22	2.33	0.43
67:O1:8:VAL:HG22	67:O1:77:ARG:HH21	2.92	0.43
68:O2:103:LYS:O	68:O2:106:VAL:HG22	5.12	0.43
69:O3:12:LYS:HD3	69:O3:96:ALA:O	2.18	0.43
71:O5:21:LEU:HD11	71:O5:55:LEU:HG	2.00	0.43
71:O5:94:LYS:O	71:O5:98:SER:N	3.53	0.43
36:1:2139:A:H62	73:O7:4:GLY:HA3	1.84	0.43
2:S0:163:ASN:ND2	2:S0:165:ARG:HG2	2.34	0.43
3:S1:103:MET:HB3	3:S1:215:VAL:HG13	2.01	0.43
6:S4:104:ASP:HB2	6:S4:108:ARG:H	4.49	0.43
6:S4:121:TYR:CG	6:S4:161:LYS:HE3	2.53	0.43
6:S4:15:PRO:HG2	6:S4:18:TRP:CE2	2.52	0.43
10:S8:103:GLN:HB3	10:S8:164:ARG:HB3	2.00	0.43
11:S9:133:HIS:O	11:S9:134:ILE:HG12	2.18	0.43
36:1:1748:G:OP2	74:O8:42:LYS:NZ	2.42	0.43
36:1:213:A:O4'	62:N6:2:ALA:HB1	2.19	0.43
36:1:2249:G:C2	36:1:2250:G:C8	3.06	0.43
36:1:2507:C:H2'	36:1:2508:U:C6	2.54	0.43
36:1:2746:A:H2'	36:1:2747:A:O4'	2.18	0.43
36:1:2984:C:H2'	36:1:2985:C:H6	1.83	0.43
36:1:95:A:C5	36:1:96:G:H1'	2.53	0.43
1:2:1174:C:C4	1:2:1175:U:C4	3.07	0.43
1:2:1180:C:H1'	1:2:1460:A:H61	1.83	0.43
1:2:1335:U:H1'	31:D9:56:ARG:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1368:G:C6	1:2:1369:U:C4	3.07	0.43
1:2:1482:C:OP2	1:2:1521:G:N1	2.44	0.43
1:2:1488:G:C8	1:2:1515:A:N6	2.87	0.43
1:2:1504:G:C2	1:2:1505:A:C4	3.06	0.43
1:2:1737:G:H2'	1:2:1738:U:O4'	2.19	0.43
1:2:489:C:H2'	1:2:490:C:C6	2.54	0.43
1:2:534:A:H5'	1:2:535:A:OP2	2.19	0.43
37:3:67:G:O5'	37:3:67:G:H8	2.02	0.43
38:4:154:C:H5''	45:L8:181:LYS:HD2	2.01	0.43
36:5:953:G:H2'	36:5:1117:G:H5''	2.00	0.43
36:5:1152:G:N2	36:5:1200:A:N6	2.64	0.43
36:5:1238:C:O2'	36:5:1239:C:OP1	2.30	0.43
36:5:1562:C:H42	36:5:1577:G:H1	1.65	0.43
36:5:1842:A:H4'	36:5:1843:C:OP2	2.19	0.43
36:5:183:G:C2	36:5:184:U:C2	3.07	0.43
39:L2:7:ASN:O	36:5:2163:C:H4'	185.95	0.43
36:5:279:U:H2'	36:5:280:U:C6	2.54	0.43
36:5:2869:U:H1'	36:5:2873:U:H5	1.83	0.43
36:5:3347:A:H61	36:5:3358:U:H3	1.67	0.43
36:5:83:U:H2'	36:5:84:U:O4'	2.18	0.43
1:6:1235:C:OP2	1:6:1245:G:H8	2.01	0.43
1:6:1268:G:H1'	1:6:1448:G:H5''	2.01	0.43
1:6:1663:G:C6	1:6:1664:C:C4	3.07	0.43
1:6:301:A:H2'	1:6:302:U:O4'	2.19	0.43
1:6:473:A:H5'	1:6:769:A:H1'	1.99	0.43
26:D4:10:ARG:HD3	1:6:780:A:N3	431.48	0.43
36:5:1055:A:H5''	37:7:100:C:O2'	2.18	0.43
38:8:141:C:H2'	38:8:142:C:C6	2.53	0.43
14:C2:60:VAL:HG22	14:C2:122:VAL:HG22	2.00	0.43
16:C4:135:ARG:HG3	1:6:1007:C:H5''	296.54	0.43
19:C7:87:GLU:O	19:C7:88:VAL:HB	3.16	0.43
23:D1:30:ALA:O	23:D1:60:ARG:NH1	4.93	0.43
4:S2:144:TRP:NE1	24:D2:97:ARG:HD2	2.33	0.43
28:D6:75:VAL:O	28:D6:79:ILE:N	2.52	0.43
28:D6:93:LYS:HB3	28:D6:93:LYS:HE2	1.68	0.43
28:D6:87:ARG:NH2	28:D6:94:ASN:HB3	3.60	0.43
41:L4:362:ASP:O	56:N0:28:ARG:HD3	3.70	0.43
41:L4:93:MET:H	41:L4:93:MET:CE	2.31	0.43
42:L5:126:GLU:HA	42:L5:196:ARG:HD2	2.01	0.43
45:L8:158:ASP:HB3	45:L8:159:PRO:HD3	2.01	0.43
46:L9:89:LYS:NZ	46:L9:191:LEU:HD12	15.68	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:M0:77:THR:OG1	47:M0:78:THR:N	2.51	0.43
48:M1:152:HIS:O	48:M1:153:LYS:HG2	5.02	0.43
48:M1:19:LEU:HD13	48:M1:40:LEU:HD12	2.00	0.43
49:M3:49:ARG:HB2	49:M3:49:ARG:HE	1.63	0.43
50:M4:8:LYS:HB3	50:M4:9:ALA:H	1.47	0.43
51:M5:68:ARG:HG2	51:M5:128:LYS:HG3	2.00	0.43
55:M9:123:LEU:O	55:M9:127:SER:N	2.47	0.43
36:1:1722:U:O4'	55:M9:96:ILE:HG23	2.18	0.43
56:N0:85:SER:OG	36:5:1294:A:H5''	301.39	0.43
60:N4:47:ARG:HG3	60:N4:54:LEU:HD23	5.22	0.43
61:N5:139:ILE:HD11	71:O5:33:VAL:HG21	1.99	0.43
45:L8:53:PRO:HD3	61:N5:32:PHE:CG	2.61	0.43
63:N7:89:VAL:HG23	63:N7:92:PHE:HE2	2.93	0.43
68:O2:102:ALA:HA	68:O2:125:ARG:HB3	3.19	0.43
2:S0:144:ILE:HG23	2:S0:158:VAL:HG13	2.48	0.43
4:S2:83:ILE:HA	4:S2:99:LYS:O	2.27	0.43
6:S4:192:ILE:HG13	6:S4:243:GLY:HA3	2.00	0.43
7:S5:99:MET:O	7:S5:100:ASN:HB2	2.42	0.43
7:S5:126:ASP:HB3	7:S5:127:GLN:H	1.87	0.43
7:S5:113:ILE:HG21	7:S5:190:ILE:HB	2.00	0.43
9:S7:16:LEU:HD22	9:S7:46:ILE:HG21	2.01	0.43
10:S8:26:LYS:O	10:S8:29:LEU:HD22	3.52	0.43
11:S9:141:VAL:HG11	11:S9:146:PHE:CG	3.20	0.43
11:S9:31:ALA:HA	11:S9:36:LEU:HB2	2.59	0.43
34:SR:40:LYS:HD2	34:SR:65:SER:HA	2.01	0.43
36:1:1341:U:H2'	36:1:1342:C:C6	2.54	0.43
36:1:241:G:C6	36:1:242:C:C4	3.07	0.43
36:1:278:U:H2'	36:1:279:U:C6	2.54	0.43
36:1:3037:U:H5''	40:L3:348:ARG:HH12	1.84	0.43
36:1:3174:A:H2'	36:1:3175:U:O4'	2.19	0.43
36:1:516:A:H2'	36:1:517:G:C8	2.54	0.43
36:1:703:G:C6	36:1:704:U:C4	3.07	0.43
1:2:1026:A:N3	1:2:1790:A:H1'	2.33	0.43
1:2:1040:G:N2	1:2:1078:C:O2	2.51	0.43
1:2:1110:G:N2	1:2:1136:U:H1'	2.34	0.43
1:2:221:A:H3'	1:2:833:U:O2	2.18	0.43
1:2:575:C:H41	25:D3:65:ASN:ND2	2.17	0.43
1:2:702:G:O2'	1:2:703:G:H8	2.02	0.43
1:2:78:A:H1'	8:S6:175:ILE:HG12	2.00	0.43
1:2:861:U:OP1	15:C3:64:ARG:NH2	2.45	0.43
1:2:870:C:H2'	1:2:871:G:C8	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:986:G:H2'	1:2:987:G:O4'	2.18	0.43
38:4:43:A:H2'	38:4:44:A:H8	1.83	0.43
51:M5:4:TYR:OH	36:5:148:G:OP2	110.24	0.43
36:5:2425:G:H2'	36:5:2426:U:O4'	2.19	0.43
36:5:174:C:N4	36:5:244:G:H1	2.15	0.43
36:5:2573:G:H5'	36:5:2574:G:OP2	2.18	0.43
36:5:283:G:O6	36:5:304:G:H1'	2.19	0.43
36:5:3353:G:H4'	36:5:3354:U:OP1	2.19	0.43
36:5:685:G:N2	36:5:696:C:C2	2.87	0.43
36:5:761:A:N1	36:5:771:A:H1'	2.34	0.43
1:6:1139:A:C5	1:6:1140:G:C8	3.07	0.43
1:6:1446:A:O2'	1:6:1447:C:H5''	2.19	0.43
1:6:1606:C:H2'	1:6:1607:G:C8	2.54	0.43
1:6:538:A:O2'	1:6:539:G:N7	2.50	0.43
1:6:654:C:H2'	1:6:655:G:H8	1.81	0.43
18:C6:10:PHE:CE2	1:6:1379:C:H5'	431.65	0.43
1:2:1605:G:OP2	18:C6:127:LYS:HE3	2.18	0.43
19:C7:6:THR:HG23	19:C7:9:VAL:HG23	2.18	0.43
20:C8:14:ILE:H	20:C8:24:GLY:N	2.17	0.43
25:D3:108:GLY:O	25:D3:109:ARG:HG2	2.18	0.43
25:D3:116:ASP:OD2	1:6:570:A:N6	358.59	0.43
26:D4:104:SER:O	26:D4:108:ARG:HG3	2.79	0.43
39:L2:107:VAL:O	39:L2:139:HIS:NE2	2.52	0.43
40:L3:56:ILE:HD12	40:L3:56:ILE:HA	2.51	0.43
37:3:121:U:H3	42:L5:268:GLU:HB3	1.82	0.43
44:L7:118:LYS:HG3	44:L7:191:VAL:CG1	2.49	0.43
36:1:2585:G:H8	45:L8:48:ARG:HG3	1.84	0.43
47:M0:75:TYR:CE2	47:M0:79:VAL:HG21	3.12	0.43
48:M1:86:VAL:HG22	48:M1:111:ASP:HB2	2.01	0.43
49:M3:149:GLN:HA	49:M3:150:PRO:HD3	1.72	0.43
51:M5:172:ARG:HH22	36:5:63:A:P	101.49	0.43
36:1:784:A:C8	54:M8:69:ARG:CZ	3.02	0.43
55:M9:103:ARG:NH2	55:M9:128:LYS:HG3	6.09	0.43
56:N0:67:ALA:O	56:N0:69:PRO:HD3	2.91	0.43
59:N3:120:LYS:HB2	59:N3:137:VAL:HG23	2.01	0.43
59:N3:27:ASP:OD2	59:N3:27:ASP:N	2.52	0.43
61:N5:40:LEU:HB3	61:N5:41:ALA:H	3.50	0.43
66:O0:22:LYS:O	66:O0:93:LEU:HB2	3.28	0.43
67:O1:82:GLU:C	67:O1:84:ASP:H	2.22	0.43
78:Q2:8:ARG:O	78:Q2:23:HIS:N	2.47	0.43
7:S5:94:THR:O	7:S5:97:LEU:HB2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:810:G:C5	9:S7:111:LYS:HE3	2.53	0.43
11:S9:131:GLN:O	11:S9:132:ARG:HG2	4.03	0.43
11:S9:168:ARG:NH2	11:S9:174:ARG:HD3	7.87	0.43
11:S9:45:ILE:HA	11:S9:45:ILE:HD13	1.81	0.43
35:SM:52:PRO:O	35:SM:54:PRO:HD3	5.05	0.43
5:S3:223:LYS:HD3	34:SR:193:ILE:HD12	2.01	0.43
36:1:1019:G:H2'	36:1:1020:G:O4'	2.19	0.43
36:1:2514:U:OP1	45:L8:68:ARG:NH1	2.47	0.43
36:1:2534:G:N2	36:1:2545:C:H42	2.15	0.43
36:1:2936:A:H2'	36:1:2937:G:C8	2.54	0.43
36:1:3082:C:H2'	36:1:3083:G:C8	2.54	0.43
36:1:3375:A:N7	36:1:3376:A:N6	2.66	0.43
1:2:1483:A:N3	1:2:1607:G:O2'	2.35	0.43
1:2:52:U:H2'	1:2:53:G:C8	2.54	0.43
37:3:105:C:O5'	37:3:105:C:H6	2.01	0.43
41:L4:183:LYS:HE3	36:5:1386:A:N7	119.68	0.43
36:5:2267:C:H2'	36:5:2268:U:C6	2.50	0.43
36:5:210:U:C2	36:5:230:U:H4'	2.54	0.43
36:5:3287:U:C2'	36:5:3288:G:H5'	2.48	0.43
36:5:3320:A:N6	36:5:3386:G:O6	2.52	0.43
36:5:599:C:H2'	36:5:600:G:O4'	2.19	0.43
36:5:594:U:H2'	36:5:609:G:O6	2.19	0.43
41:L4:112:LYS:O	36:5:790:U:H4'	122.44	0.43
36:5:815:G:C6	36:5:906:A:C4	3.06	0.43
1:6:1045:C:C2	1:6:1074:G:C2	3.07	0.43
1:6:1080:U:H2'	1:6:1081:A:C8	2.54	0.43
1:6:774:A:C5	1:6:775:G:H1'	2.54	0.43
37:7:95:A:C2	37:7:96:U:C2	3.06	0.43
13:C1:13:PHE:HE2	13:C1:15:LYS:HB3	1.83	0.43
15:C3:118:ILE:O	15:C3:122:ILE:HG13	2.19	0.43
19:C7:10:LYS:HG2	19:C7:53:TYR:CE1	2.91	0.43
22:D0:51:VAL:HG21	22:D0:94:GLU:HG3	3.50	0.43
24:D2:104:LEU:HB2	24:D2:124:LYS:O	2.18	0.43
26:D4:36:SER:O	26:D4:40:LEU:HB2	3.66	0.43
26:D4:66:GLY:HA2	1:6:532:U:H4'	431.33	0.43
27:D5:48:ASP:HA	27:D5:51:LEU:HB2	1.99	0.43
30:D8:10:ALA:HB1	30:D8:30:VAL:HB	2.03	0.43
36:1:2941:A:OP1	40:L3:255:TRP:HB3	2.18	0.43
40:L3:386:ASP:HB3	40:L3:387:LEU:H	1.56	0.43
41:L4:162:THR:O	41:L4:166:VAL:HG23	2.18	0.43
41:L4:288:ARG:O	41:L4:291:ASN:N	3.04	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:L5:111:GLN:HA	42:L5:116:ASP:HB2	2.01	0.43
42:L5:140:ARG:HH21	36:5:1080:A:P	230.18	0.43
45:L8:152:LEU:HD12	45:L8:198:ALA:HB3	2.00	0.43
45:L8:213:LYS:O	45:L8:217:THR:HG22	5.65	0.43
46:L9:93:VAL:O	46:L9:177:ASP:HA	2.18	0.43
49:M3:105:ASN:ND2	49:M3:108:ILE:HG12	2.76	0.43
49:M3:115:ARG:NH1	49:M3:147:ILE:HG12	2.34	0.43
49:M3:91:ARG:NH1	49:M3:97:VAL:HB	2.33	0.43
50:M4:103:ILE:O	50:M4:107:GLU:HB2	3.51	0.43
51:M5:27:VAL:HB	51:M5:122:ASN:HD21	1.83	0.43
52:M6:53:LYS:O	52:M6:56:ASP:HB2	2.18	0.43
59:N3:32:ARG:HE	59:N3:64:LYS:HB3	1.83	0.43
61:N5:132:ALA:O	61:N5:135:ILE:HD13	5.85	0.43
36:1:802:C:H5	64:N8:25:HIS:CD2	2.37	0.43
69:O3:53:TYR:CZ	69:O3:65:ARG:HB2	2.74	0.43
72:O6:11:LEU:HA	72:O6:11:LEU:HD13	1.77	0.43
2:S0:29:VAL:HG22	2:S0:150:ASP:HB3	1.99	0.43
5:S3:202:LEU:C	5:S3:204:ASP:H	2.36	0.43
7:S5:107:LYS:O	7:S5:111:VAL:HG23	2.18	0.43
9:S7:39:ARG:HH22	55:M9:185:LEU:HD22	1.84	0.43
11:S9:41:GLU:HG2	11:S9:44:ARG:NH2	2.70	0.43
34:SR:40:LYS:HA	34:SR:66:HIS:O	2.60	0.43
36:1:1010:G:H8	36:1:1010:G:OP2	2.02	0.43
36:1:1153:A:C6	36:1:1154:A:C6	3.07	0.43
36:1:1774:C:H2'	36:1:1775:G:O4'	2.19	0.43
36:1:1846:C:O2'	53:M7:128:ARG:NH1	2.52	0.43
36:1:2623:G:H1	36:1:2644:C:H42	1.66	0.43
36:1:871:U:H2'	36:1:872:U:C6	2.54	0.43
1:2:1209:C:H2'	1:2:1210:C:C6	2.54	0.43
1:2:1237:G:H2'	1:2:1238:A:C8	2.54	0.43
1:2:1326:A:H2'	1:2:1327:C:H6	1.83	0.43
1:2:1586:A:H2'	1:2:1587:A:O4'	2.19	0.43
1:2:246:G:O4'	6:S4:202:ASP:HB3	2.18	0.43
1:2:784:C:H2'	1:2:785:U:O4'	2.19	0.43
1:2:819:G:N2	1:2:854:U:O4'	2.51	0.43
1:2:98:U:H1'	1:2:425:A:H1'	2.01	0.43
38:4:43:A:H2'	38:4:44:A:C8	2.54	0.43
46:L9:63:LYS:NZ	36:5:1209:G:O3'	313.23	0.43
36:5:1544:G:N2	36:5:1550:C:H1'	2.34	0.43
36:5:168:U:H2'	36:5:169:U:C6	2.54	0.43
36:5:2188:A:H2'	36:5:2189:U:O4'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:2534:G:H1	36:5:2545:C:N4	2.08	0.43
36:5:115:A:C6	36:5:265:A:C6	3.07	0.43
40:L3:250:ALA:CB	36:5:2880:U:H1'	222.87	0.43
36:5:591:G:N2	36:5:612:U:OP1	2.48	0.43
1:6:1046:G:H1	1:6:1072:C:N4	2.16	0.43
1:6:1584:G:O2'	1:6:1610:G:O6	2.33	0.43
1:6:1754:A:H5'	1:6:1755:A:C8	2.53	0.43
1:6:212:U:H2'	1:6:213:A:H8	1.84	0.43
1:6:622:A:H4'	1:6:623:A:C5'	2.49	0.43
13:C1:123:VAL:HG23	13:C1:142:VAL:HA	2.01	0.43
13:C1:5:LEU:HD23	13:C1:7:VAL:H	6.17	0.43
14:C2:85:LYS:HE2	14:C2:87:PRO:HG3	4.27	0.43
15:C3:91:LEU:HA	15:C3:91:LEU:HD23	2.12	0.43
18:C6:25:GLY:H	18:C6:63:ILE:HA	1.84	0.43
25:D3:133:LEU:HD11	25:D3:137:LYS:NZ	2.34	0.43
26:D4:83:LYS:HB2	26:D4:83:LYS:HE2	4.57	0.43
6:S4:59:ARG:NH1	26:D4:87:PRO:HG3	2.32	0.43
28:D6:10:ARG:HH12	28:D6:36:ILE:HG13	3.98	0.43
28:D6:44:ILE:H	28:D6:44:ILE:HG13	1.73	0.43
29:D7:6:ASP:OD1	29:D7:9:HIS:ND1	3.02	0.43
22:D0:84:MET:HE3	31:D9:51:GLY:HA3	2.01	0.43
32:E0:14:VAL:HA	32:E0:17:GLN:HG2	2.09	0.43
39:L2:121:GLY:O	39:L2:123:ARG:HG3	2.66	0.43
39:L2:137:ILE:HG13	39:L2:147:ARG:HG2	2.01	0.43
39:L2:140:ASN:ND2	39:L2:142:ASP:HB3	6.46	0.43
39:L2:193:ARG:O	39:L2:195:SER:N	2.76	0.43
40:L3:123:TYR:CE2	40:L3:124:LYS:HG3	2.54	0.43
40:L3:218:ILE:HD11	40:L3:339:ARG:HD3	3.18	0.43
40:L3:255:TRP:HB3	36:5:2941:A:OP1	225.47	0.43
40:L3:306:THR:OG1	40:L3:316:GLU:O	2.24	0.43
41:L4:92:ASN:OD1	41:L4:92:ASN:N	3.09	0.43
45:L8:181:LYS:HG2	38:8:154:C:H5''	148.14	0.43
47:M0:10:ARG:HG2	47:M0:11:TYR:CD1	2.53	0.43
53:M7:32:THR:HG21	53:M7:87:SER:HB3	2.74	0.43
54:M8:67:ILE:HD13	54:M8:81:VAL:HG21	3.00	0.43
56:N0:53:LYS:HE3	56:N0:53:LYS:HB2	4.19	0.43
62:N6:125:LYS:O	62:N6:126:LEU:HG	2.18	0.43
38:4:23:U:H1'	62:N6:17:LYS:HE3	2.01	0.43
62:N6:52:ARG:NH2	38:8:71:A:O2'	34.52	0.43
62:N6:82:VAL:HG12	62:N6:83:ASP:O	2.19	0.43
63:N7:128:GLN:NE2	63:N7:129:TRP:H	4.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
63:N7:24:VAL:HG22	63:N7:130:PHE:CZ	2.54	0.43
65:N9:50:THR:CG2	36:5:1073:U:H1'	205.07	0.43
66:O0:43:ILE:HD12	66:O0:90:VAL:HB	1.99	0.43
67:O1:10:ARG:HH12	67:O1:44:MET:CG	6.48	0.43
68:O2:33:ARG:NH1	36:5:944:C:H4'	161.84	0.43
69:O3:45:LEU:HD23	69:O3:71:VAL:CG1	2.49	0.43
70:O4:41:ARG:NH2	70:O4:50:ALA:O	3.81	0.43
77:Q1:16:LYS:O	77:Q1:20:VAL:HG23	2.23	0.43
2:S0:57:LEU:HD12	2:S0:173:ILE:HG23	2.00	0.43
2:S0:9:LEU:HD23	2:S0:54:TRP:CD2	2.54	0.43
3:S1:183:GLN:HG2	3:S1:187:LYS:HE3	2.00	0.43
4:S2:97:ARG:HG2	4:S2:97:ARG:H	1.52	0.43
5:S3:177:MET:SD	5:S3:182:LEU:HD11	2.59	0.43
6:S4:45:ILE:HA	6:S4:61:VAL:HG11	2.00	0.43
9:S7:96:ARG:HE	9:S7:124:LYS:HB3	1.84	0.43
11:S9:129:ILE:HG12	11:S9:134:ILE:HD13	1.99	0.43
34:SR:103:PHE:CD1	34:SR:138:GLY:HA2	3.19	0.43
34:SR:294:TRP:CZ3	34:SR:301:LEU:HB2	3.01	0.43
34:SR:40:LYS:HA	34:SR:68:VAL:HG23	2.01	0.43
34:SR:90:ARG:HG2	34:SR:102:ARG:HG2	6.30	0.43
35:SM:48:ARG:NH2	36:1:1016:C:O2	2.52	0.42
36:1:1334:U:H2'	36:1:1335:C:C6	2.54	0.42
36:1:1448:U:OP1	53:M7:82:ARG:NH2	2.42	0.42
36:1:1632:A:C8	36:1:1644:C:H2'	2.54	0.42
36:1:199:A:C4	36:1:201:A:C8	3.06	0.42
36:1:207:U:H2'	36:1:208:C:H6	1.84	0.42
36:1:2536:A:H2'	36:1:2537:U:C5	2.54	0.42
36:1:3259:U:H5''	36:1:3261:C:H5	1.84	0.42
36:1:371:G:H4'	36:1:396:A:N1	2.34	0.42
36:1:573:C:C4	36:1:574:U:C4	3.06	0.42
36:1:926:A:H2'	36:1:927:C:C6	2.54	0.42
36:1:978:G:H21	36:1:979:U:H3	1.66	0.42
36:1:978:G:O2'	36:1:979:U:O2	2.35	0.42
36:1:99:A:OP1	51:M5:194:GLN:NE2	2.51	0.42
1:2:1038:U:C2	1:2:1094:G:N2	2.87	0.42
1:2:1068:C:P	3:S1:150:VAL:HG21	2.59	0.42
1:2:36:C:H5''	1:2:530:C:H5''	2.01	0.42
1:2:479:C:OP1	11:S9:121:SER:OG	2.28	0.42
36:5:1213:G:C2	36:5:1293:U:C2	3.06	0.42
36:5:1562:C:H2'	36:5:1563:C:C6	2.54	0.42
36:5:1801:U:H2'	36:5:1802:C:C6	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:2213:A:H2	36:5:2601:A:N3	2.17	0.42
36:5:2694:A:C6	36:5:2695:A:C6	3.06	0.42
36:5:595:G:H1	36:5:609:G:H5''	1.83	0.42
36:5:686:G:C2	36:5:687:U:H1'	2.53	0.42
54:M8:141:ARG:HD3	36:5:743:C:O2	174.35	0.42
36:5:96:G:H2'	36:5:97:U:O4'	2.19	0.42
1:6:320:U:H2'	1:6:321:C:C2	2.53	0.42
12:C0:3:MET:HA	12:C0:4:PRO:HD2	2.38	0.42
13:C1:46:LYS:HA	13:C1:46:LYS:HD2	2.47	0.42
15:C3:22:ALA:HB1	15:C3:23:PRO:CA	2.49	0.42
21:C9:14:PHE:CZ	21:C9:132:LEU:HD22	4.56	0.42
21:C9:33:TYR:O	21:C9:36:ILE:HG12	2.19	0.42
2:S0:185:ARG:HE	23:D1:47:PRO:HG3	1.83	0.42
24:D2:50:PHE:HB3	24:D2:63:VAL:HG22	2.00	0.42
25:D3:79:ASN:HD22	25:D3:79:ASN:N	2.17	0.42
25:D3:81:LYS:HB3	25:D3:81:LYS:HE2	1.87	0.42
28:D6:53:LEU:O	28:D6:57:SER:N	2.52	0.42
28:D6:82:ARG:HB3	28:D6:83:ILE:H	1.70	0.42
41:L4:38:VAL:HG11	41:L4:118:LYS:HA	2.01	0.42
41:L4:355:PHE:HD2	41:L4:356:THR:HG23	5.08	0.42
42:L5:183:TRP:HD1	42:L5:190:ILE:HB	6.89	0.42
42:L5:271:LYS:HD3	42:L5:271:LYS:HA	4.34	0.42
45:L8:78:PHE:C	45:L8:80:TYR:N	2.72	0.42
46:L9:103:ILE:HD11	46:L9:136:PHE:HE2	1.84	0.42
47:M0:205:SER:O	47:M0:209:ASN:HB2	2.19	0.42
49:M3:64:LYS:HG3	64:N8:69:TRP:CD2	2.54	0.42
51:M5:194:GLN:H	51:M5:194:GLN:HG2	1.58	0.42
57:N1:104:GLU:HG3	57:N1:105:PHE:N	2.33	0.42
58:N2:41:ILE:HG23	58:N2:71:PHE:HE2	3.28	0.42
60:N4:60:LYS:HB3	60:N4:63:ILE:HD11	6.93	0.42
62:N6:66:GLN:HG2	62:N6:85:VAL:HG22	5.17	0.42
64:N8:90:TYR:HB3	64:N8:100:PRO:HG3	2.01	0.42
49:M3:166:ALA:HB3	64:N8:135:GLU:HG3	2.01	0.42
68:O2:100:ILE:HG22	68:O2:105:ARG:HG2	2.00	0.42
70:O4:20:ILE:HD11	70:O4:34:HIS:NE2	3.32	0.42
1:2:1437:U:H5'	5:S3:176:LEU:HD23	2.01	0.42
7:S5:64:VAL:HG22	7:S5:89:ILE:HD11	2.00	0.42
8:S6:142:ARG:HA	8:S6:147:LEU:HD12	2.17	0.42
8:S6:43:ASP:O	8:S6:46:LYS:HB2	2.19	0.42
9:S7:11:GLN:OE1	9:S7:13:PRO:HD2	5.38	0.42
11:S9:3:ARG:HD3	11:S9:3:ARG:H	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:S9:51:LYS:HE3	11:S9:51:LYS:HB2	2.92	0.42
11:S9:4:ALA:HA	11:S9:5:PRO:HD3	1.85	0.42
34:SR:214:ALA:HB2	34:SR:220:ILE:HA	2.01	0.42
36:1:1049:C:H2'	36:1:1050:U:H6	1.84	0.42
36:1:155:G:H5''	36:1:156:G:C8	2.55	0.42
36:1:1618:G:H4'	38:4:129:C:H1'	2.00	0.42
36:1:1951:C:N4	36:1:2095:G:H1	2.17	0.42
36:1:2227:C:OP1	78:Q2:32:LYS:NZ	2.44	0.42
36:1:240:U:H4'	36:1:241:G:OP1	2.19	0.42
36:1:2514:U:OP2	36:1:2586:G:N2	2.51	0.42
36:1:2890:A:O2'	36:1:2933:A:N3	2.35	0.42
1:2:1080:U:H2'	1:2:1081:A:C8	2.53	0.42
1:2:132:U:H4'	1:2:133:U:H5'	2.01	0.42
1:2:1346:A:OP2	1:2:1348:A:N6	2.45	0.42
1:2:1443:U:O3'	1:2:1444:A:H8	2.01	0.42
1:2:144:U:H5	8:S6:137:ARG:HH11	1.66	0.42
1:2:191:C:O2'	1:2:192:U:O5'	2.34	0.42
1:2:492:A:H4'	1:2:493:U:C5	2.54	0.42
1:2:551:G:H1	1:2:573:C:N4	2.13	0.42
1:2:67:A:N6	1:2:83:G:O2'	2.52	0.42
1:2:43:A:O2'	1:2:99:C:OP1	2.29	0.42
36:5:1028:U:H2'	36:5:1028:U:O2	2.19	0.42
36:5:951:A:H5''	36:5:1143:A:N1	2.34	0.42
36:5:1155:C:H2'	36:5:1156:C:C6	2.54	0.42
36:5:1159:A:O2'	36:5:1160:C:H5'	2.19	0.42
36:5:1566:A:C2'	36:5:1567:U:H5'	2.49	0.42
36:5:3280:U:O2'	36:5:3281:U:H6	2.01	0.42
36:5:379:C:H2'	36:5:380:U:C6	2.51	0.42
36:5:565:U:C2	36:5:566:G:C8	3.07	0.42
64:N8:111:LYS:HE2	36:5:714:G:N7	139.57	0.42
36:5:835:G:O2'	36:5:857:G:N2	2.29	0.42
54:M8:144:ARG:HH12	36:5:976:U:H5'	177.65	0.42
3:S1:146:GLN:NE2	1:6:1066:C:O4'	343.07	0.42
1:6:1076:A:H2'	1:6:1077:C:O4'	2.19	0.42
1:6:1360:A:H2'	1:6:1361:U:H4'	2.00	0.42
31:D9:55:PHE:O	1:6:1418:G:N2	413.35	0.42
1:6:1433:G:H2'	1:6:1434:U:O4'	2.20	0.42
7:S5:102:ARG:HH12	1:6:1474:G:H5'	353.58	0.42
1:6:1684:U:H2'	1:6:1685:G:C8	2.54	0.42
1:6:1757:G:H2'	1:6:1758:U:H6	1.83	0.42
1:6:485:A:H61	1:6:502:U:H3	1.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:922:G:H2'	1:6:923:A:H8	1.83	0.42
37:7:94:C:H2'	37:7:95:A:H8	1.85	0.42
12:C0:44:LYS:HE3	1:6:1217:A:H4'	425.57	0.42
14:C2:79:ALA:HB1	14:C2:87:PRO:O	2.19	0.42
15:C3:42:ARG:HB2	15:C3:42:ARG:HH11	1.84	0.42
19:C7:48:ASN:HD22	1:6:1388:A:H5''	429.87	0.42
22:D0:26:LEU:HB3	22:D0:34:LEU:HD21	2.00	0.42
26:D4:122:GLY:HA2	26:D4:125:LEU:HB2	3.80	0.42
27:D5:59:TYR:HE2	27:D5:100:ILE:HA	1.84	0.42
28:D6:10:ARG:HH22	28:D6:35:ALA:C	6.98	0.42
1:2:1418:G:O2'	31:D9:56:ARG:O	2.35	0.42
36:1:2424:A:N1	39:L2:230:VAL:HG21	2.35	0.42
40:L3:377:HIS:NE2	40:L3:387:LEU:HD11	2.34	0.42
41:L4:194:TYR:N	41:L4:194:TYR:CD2	2.87	0.42
41:L4:309:ARG:HH21	41:L4:312:VAL:CB	3.41	0.42
41:L4:317:PRO:HA	41:L4:323:VAL:HG13	2.22	0.42
43:L6:175:LYS:HA	43:L6:175:LYS:HD2	4.53	0.42
43:L6:175:LYS:HE3	43:L6:175:LYS:HB3	1.79	0.42
43:L6:49:GLY:O	43:L6:163:PHE:N	2.59	0.42
49:M3:47:ALA:HB1	49:M3:48:PRO:CD	2.48	0.42
51:M5:168:GLY:O	51:M5:172:ARG:HB2	2.47	0.42
52:M6:80:PHE:HD2	52:M6:104:VAL:HG11	1.84	0.42
54:M8:122:ILE:HG12	54:M8:123:THR:O	5.39	0.42
36:1:1940:G:OP1	55:M9:80:LYS:HE3	2.19	0.42
56:N0:47:LYS:O	56:N0:48:LEU:HD23	2.18	0.42
59:N3:26:ALA:O	59:N3:115:THR:HG23	2.19	0.42
66:O0:23:TYR:HB2	66:O0:91:SER:O	2.19	0.42
67:O1:20:LEU:O	67:O1:22:GLY:N	2.52	0.42
67:O1:70:ARG:HE	67:O1:102:LYS:HZ3	6.01	0.42
67:O1:88:PRO:HG2	67:O1:89:LEU:HD13	2.00	0.42
73:O7:14:LYS:H	73:O7:14:LYS:HG2	1.45	0.42
74:O8:12:LEU:HA	74:O8:12:LEU:HD13	2.15	0.42
2:S0:193:GLN:O	2:S0:195:TRP:N	2.52	0.42
2:S0:8:ASP:HB3	2:S0:9:LEU:H	2.26	0.42
3:S1:109:LYS:O	3:S1:113:MET:HG3	2.19	0.42
5:S3:73:VAL:HG13	5:S3:77:PHE:HD2	2.57	0.42
9:S7:71:HIS:HA	9:S7:74:GLN:HB2	3.17	0.42
11:S9:126:ARG:HG3	32:E0:33:ARG:HD3	2.54	0.42
11:S9:7:THR:OG1	11:S9:8:TYR:N	2.52	0.42
36:1:1821:U:C4	70:O4:67:LYS:HE2	2.53	0.42
36:1:2213:A:H2'	36:1:2214:A:H8	1.78	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:99:A:H1'	36:1:281:G:N7	2.34	0.42
36:1:3070:A:C6	36:1:3071:U:C4	3.08	0.42
36:1:87:U:H2'	36:1:88:A:C8	2.55	0.42
36:1:8:C:H2'	36:1:9:U:C6	2.55	0.42
1:2:1308:G:C2	1:2:1309:C:C2	3.07	0.42
1:2:268:C:H41	8:S6:186:ARG:HD3	1.84	0.42
1:2:361:C:H2'	1:2:362:G:H8	1.82	0.42
1:2:463:U:H2'	1:2:464:A:H8	1.84	0.42
1:2:503:G:O2'	1:2:504:U:OP1	2.29	0.42
38:4:29:U:H5''	49:M3:27:ASP:HB3	2.01	0.42
36:5:1579:C:H5''	36:5:1649:U:H5''	2.01	0.42
36:5:189:G:H3'	36:5:224:C:OP2	2.19	0.42
36:5:2197:C:N4	36:5:2241:U:H2'	2.35	0.42
51:M5:87:GLN:NE2	36:5:2422:C:O2'	172.30	0.42
36:5:2815:G:H5''	36:5:2816:G:OP2	2.18	0.42
36:5:2830:G:H1'	36:5:2861:U:C2	2.54	0.42
36:5:29:C:H4'	36:5:62:A:H4'	2.02	0.42
36:5:3198:U:H4'	36:5:3199:G:OP2	2.19	0.42
36:5:963:G:H8	36:5:963:G:OP2	2.02	0.42
1:6:187:G:H8	1:6:187:G:O5'	2.03	0.42
1:6:56:U:O2	1:6:57:G:H1'	2.18	0.42
9:S7:104:ARG:HG2	1:6:742:U:H1'	352.75	0.42
38:8:126:A:O2'	38:8:129:C:N4	2.52	0.42
71:O5:56:THR:HG23	38:8:60:U:N3	51.27	0.42
13:C1:53:TYR:CG	13:C1:113:PRO:HG2	2.53	0.42
16:C4:85:ALA:H	16:C4:119:THR:CG2	2.30	0.42
19:C7:35:CYS:HA	19:C7:38:ILE:HG22	2.02	0.42
22:D0:33:GLN:N	22:D0:33:GLN:OE1	2.51	0.42
24:D2:53:ILE:HB	24:D2:60:LYS:HB2	4.51	0.42
26:D4:72:PHE:HE1	26:D4:74:LEU:HD23	2.26	0.42
29:D7:49:HIS:HD2	1:6:958:U:H5'	342.21	0.42
31:D9:33:LYS:HE2	31:D9:34:TYR:CZ	4.14	0.42
32:E0:13:LYS:HB3	32:E0:17:GLN:NE2	6.59	0.42
39:L2:112:ILE:HD12	79:Q3:79:VAL:HG22	2.00	0.42
40:L3:97:ARG:HE	40:L3:97:ARG:HB2	1.69	0.42
41:L4:20:LEU:HA	41:L4:21:PRO:HD3	1.93	0.42
41:L4:22:LEU:HD22	41:L4:23:PRO:CD	2.49	0.42
45:L8:134:TYR:CD2	45:L8:134:TYR:N	2.91	0.42
36:1:2523:A:O2'	45:L8:49:TYR:O	2.31	0.42
46:L9:86:TYR:CG	46:L9:151:VAL:HG13	3.08	0.42
46:L9:89:LYS:HD2	46:L9:183:HIS:ND1	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:172:ARG:CZ	51:M5:174:ILE:HD11	2.82	0.42
53:M7:3:ARG:HD2	53:M7:3:ARG:HA	4.80	0.42
54:M8:85:GLY:O	54:M8:104:LEU:HB2	3.60	0.42
55:M9:148:ASP:OD1	55:M9:151:ARG:NH2	4.34	0.42
55:M9:77:GLY:HA3	36:5:1939:G:OP1	218.31	0.42
57:N1:45:ASN:OD1	57:N1:47:SER:OG	3.12	0.42
57:N1:91:LEU:HD12	57:N1:96:ILE:HD11	2.00	0.42
61:N5:53:HIS:CE1	61:N5:56:ARG:HG2	2.55	0.42
61:N5:67:ILE:CD1	61:N5:121:LYS:HG3	2.84	0.42
64:N8:77:LYS:HD3	64:N8:77:LYS:HA	1.87	0.42
65:N9:14:ARG:NH1	65:N9:18:ARG:HH11	2.08	0.42
66:O0:41:LEU:HD23	66:O0:66:LYS:O	3.85	0.42
71:O5:119:LYS:HZ2	71:O5:119:LYS:HA	5.02	0.42
71:O5:95:PHE:CD2	36:5:136:G:H5'	63.60	0.42
72:O6:53:TYR:CD1	72:O6:76:ARG:HG2	2.54	0.42
3:S1:106:THR:O	3:S1:109:LYS:N	2.52	0.42
3:S1:70:LEU:HA	3:S1:73:LEU:HB2	4.82	0.42
3:S1:70:LEU:HD13	3:S1:79:HIS:CD2	4.92	0.42
5:S3:46:THR:N	5:S3:83:THR:O	2.86	0.42
7:S5:166:ARG:HB2	30:D8:46:GLY:HA3	2.00	0.42
7:S5:198:LEU:HD23	7:S5:198:LEU:HA	2.13	0.42
8:S6:139:ASN:HA	8:S6:142:ARG:HB2	2.21	0.42
9:S7:16:LEU:O	9:S7:20:VAL:HG23	2.19	0.42
10:S8:167:ALA:HA	10:S8:183:ILE:HA	2.46	0.42
34:SR:23:LEU:HD13	34:SR:302:PHE:HB3	2.94	0.42
36:1:1062:A:N3	57:N1:130:ARG:NH2	2.66	0.42
36:1:1108:U:H2'	36:1:1109:U:H6	1.85	0.42
36:1:1894:U:O2'	36:1:3054:U:OP1	2.36	0.42
36:1:3216:G:O6	36:1:3259:U:H2'	2.20	0.42
36:1:597:G:C2	36:1:608:A:H1'	2.53	0.42
36:1:830:A:H2'	36:1:831:G:O4'	2.20	0.42
36:1:863:C:H2'	36:1:864:G:O4'	2.19	0.42
1:2:1183:A:C6	1:2:1184:A:N1	2.87	0.42
1:2:409:C:H2'	1:2:410:A:C8	2.55	0.42
1:2:938:G:N2	1:2:940:A:H3'	2.35	0.42
36:5:1157:G:H2'	36:5:1158:A:O4'	2.19	0.42
36:5:1644:C:H5'	36:5:1645:U:O5'	2.20	0.42
36:5:2631:U:H4'	36:5:2697:A:H2	1.84	0.42
46:L9:62:ARG:NH1	36:5:3114:A:O2'	324.85	0.42
36:5:3192:U:H2'	36:5:3193:C:C6	2.54	0.42
36:5:3269:U:H5'	36:5:3271:G:O4'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:676:G:N2	36:5:786:A:C4	2.87	0.42
1:6:151:G:N2	1:6:163:G:N2	2.67	0.42
8:S6:59:GLN:HA	1:6:155:U:H4'	304.43	0.42
1:6:1196:A:O2'	1:6:1602:C:O3'	2.37	0.42
1:6:74:U:C2	1:6:76:A:H5''	2.55	0.42
29:D7:49:HIS:CD2	1:6:958:U:H5'	342.81	0.42
42:L5:158:ARG:HD2	37:7:47:C:OP2	284.89	0.42
38:8:91:C:H2'	38:8:92:A:C8	2.54	0.42
15:C3:94:LYS:HE2	1:6:953:G:P	301.86	0.42
1:2:887:A:H5''	16:C4:120:PRO:HB2	2.00	0.42
16:C4:11:SER:HG	16:C4:12:GLN:H	4.10	0.42
18:C6:123:ARG:HG3	18:C6:124:PRO:HD2	2.00	0.42
19:C7:23:LYS:HG3	34:SR:198:ASN:HD21	1.84	0.42
21:C9:33:TYR:HB2	21:C9:37:VAL:HG23	4.80	0.42
24:D2:6:VAL:HG12	24:D2:34:ILE:HD11	2.00	0.42
28:D6:11:ASN:O	28:D6:33:ASP:HB2	2.43	0.42
30:D8:26:THR:O	30:D8:44:VAL:HG22	3.45	0.42
33:E1:109:ASP:HB2	33:E1:113:LYS:HD2	3.79	0.42
40:L3:148:LEU:HD12	40:L3:148:LEU:HA	1.95	0.42
40:L3:255:TRP:CD1	40:L3:256:HIS:CE1	3.46	0.42
40:L3:30:LYS:NZ	36:5:3139:A:OP2	235.48	0.42
41:L4:10:SER:OG	41:L4:14:GLU:O	4.12	0.42
42:L5:107:ARG:HH12	42:L5:120:LYS:HA	1.85	0.42
42:L5:130:GLU:O	42:L5:131:LEU:HD22	2.19	0.42
36:1:2688:U:N3	42:L5:17:GLN:O	2.25	0.42
42:L5:285:ARG:HH12	37:7:62:U:H4'	341.63	0.42
43:L6:154:LEU:HD23	43:L6:157:GLN:HB2	2.00	0.42
45:L8:43:LYS:HD3	45:L8:43:LYS:HA	1.78	0.42
46:L9:57:VAL:HG23	46:L9:68:LEU:HG	2.02	0.42
46:L9:69:ARG:HA	46:L9:69:ARG:HD2	1.60	0.42
47:M0:210:ILE:HG23	47:M0:217:PHE:CD2	2.54	0.42
36:1:1128:U:O4'	47:M0:4:ARG:NH2	2.53	0.42
48:M1:96:PHE:CD1	48:M1:102:PHE:HB3	2.54	0.42
50:M4:47:ASP:OD1	50:M4:78:THR:HA	2.90	0.42
52:M6:39:GLU:N	52:M6:39:GLU:OE1	2.35	0.42
59:N3:15:LEU:HD23	59:N3:53:SER:HB3	2.07	0.42
63:N7:16:GLY:O	70:O4:74:ARG:HG3	5.23	0.42
68:O2:2:ALA:HB1	68:O2:90:LYS:HE3	2.01	0.42
72:O6:26:ILE:H	72:O6:26:ILE:HG13	1.49	0.42
74:O8:65:LEU:O	74:O8:65:LEU:HD23	2.67	0.42
75:O9:48:LYS:HD3	75:O9:48:LYS:HA	3.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
79:Q3:49:ARG:HB2	79:Q3:55:TRP:CZ3	2.54	0.42
2:S0:126:PRO:HG3	2:S0:151:SER:HB3	2.99	0.42
5:S3:106:LYS:HG3	5:S3:175:VAL:HB	2.01	0.42
5:S3:211:PRO:O	5:S3:212:LYS:HD3	5.31	0.42
7:S5:102:ARG:HG3	7:S5:102:ARG:H	2.77	0.42
7:S5:183:ALA:HB3	7:S5:190:ILE:HD13	2.63	0.42
9:S7:116:ARG:HD3	9:S7:116:ARG:HA	4.53	0.42
11:S9:123:HIS:CD2	32:E0:33:ARG:HD2	4.99	0.42
11:S9:68:LYS:NZ	11:S9:72:GLU:OE1	2.45	0.42
1:2:1460:A:H1'	35:SM:71:ASN:O	2.18	0.42
36:1:1115:G:OP1	64:N8:22:ILE:HG21	2.18	0.42
36:1:1458:U:H5'	67:O1:30:PRO:HB3	2.01	0.42
36:1:1930:A:H5'	36:1:1932:A:O2'	2.19	0.42
36:1:2093:A:H2'	36:1:2094:C:O4'	2.19	0.42
36:1:225:C:H4'	62:N6:32:SER:O	2.20	0.42
36:1:241:G:C6	36:1:242:C:N4	2.88	0.42
36:1:3168:A:H3'	36:1:3169:U:C6	2.55	0.42
36:1:3253:G:O5'	36:1:3253:G:H8	2.03	0.42
36:1:373:A:N1	36:1:394:G:H4'	2.35	0.42
36:1:435:C:H2'	36:1:436:A:C8	2.55	0.42
1:2:1064:G:H2'	1:2:1065:A:C8	2.54	0.42
1:2:1111:G:C6	1:2:1112:G:C4	3.07	0.42
1:2:145:A:HO2'	1:2:146:U:P	2.39	0.42
1:2:1553:G:N1	1:2:1556:A:OP2	2.45	0.42
1:2:1559:A:N6	20:C8:134:ARG:HD2	2.34	0.42
1:2:1699:G:H2'	1:2:1700:C:H5''	2.02	0.42
1:2:137:U:OP1	1:2:1706:C:H5'	2.20	0.42
37:3:115:G:H2'	37:3:116:C:H6	1.84	0.42
36:5:1064:A:H5''	36:5:1066:G:O4'	2.19	0.42
36:5:2198:A:C8	36:5:2270:A:H1'	2.55	0.42
36:5:247:C:C4	36:5:248:U:H1'	2.54	0.42
36:5:2607:G:C4	36:5:2608:G:C8	3.08	0.42
36:5:3231:U:H2'	36:5:3232:G:C8	2.54	0.42
36:5:3281:U:H5'	36:5:3282:U:OP2	2.19	0.42
40:L3:334:ARG:NH2	36:5:3304:U:O2'	212.60	0.42
36:5:441:U:O2'	36:5:492:U:O2	2.30	0.42
1:6:1037:C:H2'	1:6:1038:U:H6	1.85	0.42
1:6:1063:U:H2'	1:6:1064:G:H8	1.85	0.42
1:6:1083:G:H1	1:6:1090:C:H42	1.67	0.42
1:6:489:C:O2'	1:6:490:C:O5'	2.37	0.42
1:6:689:G:H2'	1:6:690:G:O4'	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:643:G:H1	1:6:691:C:H42	1.66	0.42
1:6:877:G:H5'	1:6:937:C:H1'	2.00	0.42
14:C2:45:LEU:HB3	14:C2:46:ARG:H	2.76	0.42
14:C2:73:LYS:HE3	33:E1:108:VAL:HG13	2.02	0.42
16:C4:91:THR:O	16:C4:93:THR:N	2.79	0.42
22:D0:43:LYS:O	22:D0:47:GLN:HB2	2.78	0.42
22:D0:82:TYR:HB3	31:D9:52:PHE:HB3	2.08	0.42
25:D3:50:LYS:HB3	25:D3:103:LEU:HD23	4.07	0.42
25:D3:43:PHE:CE1	25:D3:49:ALA:HB3	2.71	0.42
32:E0:50:VAL:HA	32:E0:54:ARG:HA	3.72	0.42
39:L2:120:PRO:HB3	39:L2:162:ALA:HA	2.01	0.42
40:L3:37:ARG:CG	40:L3:187:SER:H	4.35	0.42
36:1:2395:G:H4'	40:L3:258:ALA:HB1	2.00	0.42
40:L3:306:THR:HA	40:L3:307:PRO:HD3	1.97	0.42
41:L4:222:VAL:HA	41:L4:223:PRO:HD3	2.07	0.42
41:L4:22:LEU:CD1	41:L4:26:PHE:HB2	2.73	0.42
42:L5:285:ARG:NH1	37:7:62:U:O3'	341.71	0.42
42:L5:48:LYS:HZ1	36:5:2749:G:P	241.67	0.42
36:1:1103:A:N9	44:L7:158:LYS:HD3	2.35	0.42
44:L7:224:ILE:HG23	56:N0:36:ILE:HG12	2.64	0.42
45:L8:134:TYR:CG	45:L8:190:VAL:HG21	2.54	0.42
46:L9:122:LYS:HG2	46:L9:123:ILE:N	2.35	0.42
47:M0:34:TYR:O	47:M0:88:ARG:HA	2.42	0.42
52:M6:105:PHE:HE2	52:M6:112:TYR:CE2	3.74	0.42
52:M6:18:ARG:NH1	52:M6:128:ARG:HD2	2.34	0.42
52:M6:25:LYS:HD3	52:M6:25:LYS:O	2.65	0.42
54:M8:26:LEU:HD23	54:M8:29:LEU:HD12	3.29	0.42
57:N1:12:ARG:HD3	57:N1:13:TYR:CZ	4.76	0.42
60:N4:45:ASN:HA	60:N4:46:PRO:HD3	2.00	0.42
69:O3:42:GLN:OE1	69:O3:45:LEU:HD12	2.18	0.42
73:O7:76:ASN:O	73:O7:79:GLN:HG3	2.48	0.42
77:Q1:14:LYS:HD2	1:6:1115:U:H5''	295.16	0.42
77:Q1:12:ARG:HG2	77:Q1:15:ARG:NH2	2.34	0.42
39:L2:177:LYS:NZ	79:Q3:26:VAL:HG13	2.34	0.42
2:S0:131:GLN:O	2:S0:135:GLU:HB2	2.69	0.42
2:S0:74:VAL:HG12	2:S0:76:ILE:HG12	2.14	0.42
3:S1:45:LYS:HB2	16:C4:13:VAL:HG23	2.01	0.42
6:S4:123:LEU:HA	6:S4:160:VAL:O	2.20	0.42
6:S4:159:THR:HG23	6:S4:173:ILE:HB	2.01	0.42
7:S5:143:ARG:NH1	7:S5:218:GLU:OE1	2.53	0.42
10:S8:155:SER:O	10:S8:159:GLN:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:S9:171:ARG:HH11	11:S9:174:ARG:HD3	5.76	0.42
36:1:1028:U:O2'	48:M1:94:ARG:NH2	2.52	0.42
36:1:1404:G:N2	36:1:1407:A:OP2	2.49	0.42
36:1:116:A:H62	36:1:153:U:H1'	1.84	0.42
36:1:1833:G:OP1	75:O9:10:LYS:NZ	2.42	0.42
36:1:3121:U:C2	36:1:3122:A:N7	2.87	0.42
36:1:3340:G:H5''	36:1:3341:U:OP1	2.19	0.42
36:1:528:U:H2'	36:1:529:A:H8	1.79	0.42
36:1:674:G:C5	36:1:675:C:C4	3.08	0.42
36:1:856:G:C6	36:1:857:G:N1	2.87	0.42
1:2:1054:U:C4	1:2:1055:U:C4	3.06	0.42
1:2:1183:A:N1	17:C5:99:GLY:HA3	2.35	0.42
1:2:1183:A:N6	1:2:1184:A:N1	2.67	0.42
1:2:1202:A:N3	1:2:1202:A:H3'	2.35	0.42
1:2:1773:C:OP1	77:Q1:3:ALA:HB3	2.20	0.42
1:2:588:U:H2'	1:2:589:C:O4'	2.20	0.42
1:2:95:G:O2'	1:2:460:A:O2'	2.35	0.42
37:3:28:C:H1'	37:3:55:A:H61	1.83	0.42
36:5:1440:G:H2'	36:5:1441:G:C8	2.55	0.42
36:5:1712:G:C2	36:5:1713:G:N2	2.88	0.42
36:5:1908:A:H2'	36:5:1909:A:O4'	2.19	0.42
36:5:1946:A:C6	36:5:1947:G:C6	3.07	0.42
36:5:3060:C:H2'	36:5:3061:G:H8	1.85	0.42
36:5:3192:U:H2'	36:5:3193:C:H6	1.85	0.42
36:5:32:U:H6	36:5:32:U:O5'	2.02	0.42
36:5:378:A:N7	36:5:391:A:H2	2.17	0.42
36:5:507:U:H2'	36:5:508:U:C6	2.55	0.42
36:5:64:G:C5	36:5:322:U:C4	3.08	0.42
36:5:734:C:H2'	36:5:735:A:O4'	2.20	0.42
33:E1:134:ASN:O	1:6:1251:U:H5''	439.23	0.42
1:6:1344:A:O2'	1:6:1345:A:OP1	2.29	0.42
1:6:1361:U:H5'	1:6:1362:U:OP2	2.19	0.42
1:6:1394:G:O2'	1:6:1395:G:H5'	2.20	0.42
1:6:1685:G:H8	1:6:1685:G:O5'	2.02	0.42
1:6:1757:G:H2'	1:6:1758:U:C6	2.55	0.42
1:6:780:A:H5''	1:6:781:U:H5'	2.02	0.42
15:C3:94:LYS:NZ	1:6:952:A:OP1	300.11	0.42
18:C6:7:VAL:N	18:C6:22:VAL:O	2.89	0.42
21:C9:84:LYS:HD2	21:C9:94:ILE:HG13	5.53	0.42
23:D1:36:VAL:HB	23:D1:51:VAL:HB	2.14	0.42
23:D1:3:ASN:HB2	23:D1:4:ASP:H	1.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:D2:18:GLU:HG2	24:D2:65:LEU:HG	2.56	0.42
22:D0:83:GLU:OE2	31:D9:55:PHE:HB2	2.18	0.42
33:E1:143:LYS:O	33:E1:145:HIS:N	2.52	0.42
39:L2:166:ILE:H	39:L2:166:ILE:HG12	1.80	0.42
40:L3:151:ILE:O	40:L3:155:ALA:HB3	2.20	0.42
40:L3:205:VAL:HG11	40:L3:322:ILE:CD1	3.17	0.42
40:L3:21:ARG:HG3	36:5:2991:A:OP1	209.87	0.42
40:L3:54:THR:OG1	40:L3:55:THR:N	2.51	0.42
42:L5:259:LYS:HE2	42:L5:259:LYS:HB3	1.89	0.42
45:L8:122:LYS:C	45:L8:124:ASP:H	2.21	0.42
46:L9:134:ILE:HG12	46:L9:146:LEU:HG	2.01	0.42
46:L9:81:GLY:O	46:L9:85:GLY:HA2	2.40	0.42
47:M0:207:GLU:HB3	47:M0:211:ARG:NH2	4.85	0.42
50:M4:118:PHE:O	50:M4:122:VAL:HG23	2.23	0.42
51:M5:114:ARG:HG2	51:M5:137:PRO:HG3	2.04	0.42
52:M6:106:GLU:HG2	52:M6:106:GLU:H	1.61	0.42
52:M6:180:SER:O	52:M6:184:THR:HG23	6.43	0.42
54:M8:34:THR:HG23	54:M8:49:LEU:HD21	2.02	0.42
36:1:523:A:O2'	56:N0:69:PRO:HD2	2.20	0.42
59:N3:66:LYS:HB2	59:N3:69:LEU:HD22	2.02	0.42
70:O4:58:ARG:O	70:O4:61:GLN:HB2	2.82	0.42
72:O6:98:ARG:HB3	72:O6:99:ARG:H	4.39	0.42
76:Q0:104:PRO:HB2	76:Q0:107:ALA:HB2	2.14	0.42
77:Q1:8:LYS:HD3	77:Q1:12:ARG:NH2	2.41	0.42
79:Q3:8:VAL:HG22	36:5:1927:G:OP1	245.22	0.42
2:S0:121:VAL:HG23	2:S0:141:ILE:HG21	2.01	0.42
5:S3:7:LYS:HA	5:S3:10:LYS:HB3	2.01	0.42
6:S4:49:ARG:HH12	1:6:448:C:P	377.66	0.42
7:S5:185:ARG:HA	1:6:1535:U:H5	335.44	0.42
34:SR:132:LYS:HB2	34:SR:134:TRP:HE1	1.83	0.42
34:SR:295:SER:HB3	34:SR:302:PHE:CE2	4.23	0.42
34:SR:36:ALA:HB1	34:SR:68:VAL:HG12	2.01	0.42
36:1:1080:A:OP1	42:L5:140:ARG:HD3	2.19	0.42
36:1:1230:G:H1	36:1:1279:C:N4	2.13	0.42
36:1:143:G:C2	36:1:144:A:C8	3.07	0.42
36:1:1580:A:H1'	36:1:1581:C:C5	2.54	0.42
36:1:153:U:O3'	36:1:158:G:H4'	2.20	0.42
36:1:184:U:H2'	36:1:185:C:H6	1.85	0.42
36:1:2390:A:H2'	36:1:2391:G:O4'	2.20	0.42
36:1:3353:G:C5	36:1:3356:G:C8	3.07	0.42
36:1:801:A:H4'	36:1:802:C:O5'	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1370:U:H4'	1:2:1371:A:O5'	2.17	0.42
1:2:1464:G:O3'	18:C6:141:SER:OG	2.27	0.42
1:2:1488:G:H5'	1:2:1489:U:OP1	2.20	0.42
1:2:831:U:O2	1:2:831:U:H2'	2.18	0.42
1:2:916:U:H3	16:C4:41:ARG:NH2	2.18	0.42
1:2:934:C:C4	1:2:1077:C:H4'	2.55	0.42
38:4:17:A:H2'	38:4:18:U:O4'	2.19	0.42
36:5:1307:G:O2'	36:5:1308:A:N7	2.48	0.42
36:5:144:A:H2'	36:5:145:G:O4'	2.19	0.42
36:5:1567:U:H1'	36:5:1570:U:O4	2.20	0.42
39:L2:70:ARG:HB3	36:5:1650:G:H5''	179.48	0.42
36:5:1688:U:H2'	36:5:1689:U:C6	2.54	0.42
36:5:1728:G:H5''	36:5:1730:G:O4'	2.19	0.42
36:5:2594:C:H2'	36:5:2595:A:O4'	2.18	0.42
57:N1:57:TYR:OH	36:5:2724:U:OP1	222.98	0.42
36:5:2794:G:O2'	36:5:2795:U:OP2	2.33	0.42
36:5:3127:A:H2'	36:5:3128:G:O4'	2.20	0.42
36:5:422:A:C2	36:5:2363:A:H4'	2.55	0.42
36:5:618:C:H2'	36:5:619:A:C8	2.54	0.42
36:5:660:A:C2	36:5:1435:A:C2	3.07	0.42
36:5:697:A:C2	36:5:698:U:C2	3.07	0.42
36:5:811:U:H2'	36:5:812:G:H8	1.83	0.42
1:6:1207:C:H42	1:6:1456:C:H5	1.66	0.42
1:6:1350:U:H2'	1:6:1351:G:C8	2.53	0.42
1:6:1451:C:H2'	1:6:1452:U:C6	2.54	0.42
1:6:1543:A:H1'	1:6:1569:A:C2	2.55	0.42
17:C5:115:TYR:OH	1:6:1556:A:OP1	388.18	0.42
1:6:217:A:O2'	1:6:218:A:H8	2.03	0.42
1:6:72:A:H2'	1:6:73:U:C6	2.54	0.42
38:8:138:A:C2	38:8:139:U:C2	3.07	0.42
38:8:5:U:H2'	38:8:6:U:C6	2.55	0.42
12:C0:23:ALA:CB	12:C0:64:TYR:HB2	3.62	0.42
12:C0:47:GLN:O	12:C0:50:THR:HG22	2.19	0.42
14:C2:54:ARG:HD3	14:C2:56:GLU:CD	4.03	0.42
1:2:960:U:O2'	15:C3:51:GLY:HA3	2.20	0.42
20:C8:108:LYS:N	20:C8:108:LYS:HD2	2.35	0.42
20:C8:81:ILE:O	20:C8:83:ALA:N	2.45	0.42
21:C9:135:ILE:HG13	21:C9:135:ILE:H	2.33	0.42
27:D5:58:ARG:HG2	27:D5:103:ARG:HH11	3.66	0.42
27:D5:55:PRO:O	27:D5:56:THR:OG1	2.34	0.42
28:D6:88:SER:OG	28:D6:91:ASP:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:D9:14:TYR:CD2	31:D9:14:TYR:N	2.87	0.42
33:E1:113:LYS:HD2	33:E1:113:LYS:H	3.21	0.42
39:L2:181:LYS:HG3	39:L2:184:ARG:HG3	2.01	0.42
40:L3:161:LEU:HD22	40:L3:178:LEU:HD11	2.02	0.42
40:L3:311:PHE:HE2	40:L3:317:ILE:HG12	3.92	0.42
36:1:3304:U:P	40:L3:332:ARG:HH22	2.42	0.42
42:L5:207:TYR:O	42:L5:211:LEU:N	2.53	0.42
44:L7:136:TYR:CE2	44:L7:231:ASN:HB2	2.54	0.42
45:L8:108:ARG:O	45:L8:112:GLU:HG2	2.20	0.42
45:L8:75:ILE:HD11	51:M5:22:LEU:HD21	2.92	0.42
36:1:1010:G:N2	47:M0:193:ASP:OD2	2.52	0.42
36:1:2350:C:O3'	53:M7:68:GLY:HA3	2.20	0.42
55:M9:127:SER:HA	55:M9:132:PHE:CD2	2.53	0.42
55:M9:6:THR:O	55:M9:10:LEU:HB2	2.20	0.42
57:N1:7:TYR:OH	57:N1:54:HIS:HB2	2.74	0.42
59:N3:2:SER:CA	59:N3:56:ASP:HA	5.64	0.42
60:N4:42:GLN:HB2	60:N4:42:GLN:HE21	3.27	0.42
62:N6:48:LEU:HG	62:N6:122:LYS:NZ	4.68	0.42
36:1:641:C:OP1	64:N8:21:ARG:HB3	2.20	0.42
36:1:1730:G:C6	66:O0:26:GLY:HA3	2.55	0.42
67:O1:11:GLU:OE2	67:O1:74:ARG:NH2	2.70	0.42
36:1:624:G:P	69:O3:86:ARG:HH22	2.43	0.42
2:S0:29:VAL:HG11	2:S0:37:VAL:HG21	12.15	0.42
3:S1:193:ILE:HG22	3:S1:197:ILE:HD13	4.85	0.42
5:S3:175:VAL:HG13	5:S3:182:LEU:HB2	2.01	0.42
5:S3:202:LEU:HA	5:S3:203:PRO:HD3	2.06	0.42
5:S3:43:PRO:O	5:S3:44:THR:HB	2.43	0.42
9:S7:44:LYS:HE3	9:S7:63:PRO:HA	2.01	0.42
9:S7:62:VAL:HG12	9:S7:64:VAL:H	1.85	0.42
11:S9:87:SER:HB3	11:S9:90:LYS:HB2	4.71	0.42
34:SR:59:ARG:NH2	34:SR:96:THR:O	2.52	0.42
36:1:1316:C:C4	52:M6:130:LYS:HA	2.55	0.42
36:1:1653:G:H2'	36:1:1654:A:C8	2.54	0.42
36:1:170:G:H1	36:1:248:U:H3	1.67	0.42
36:1:2696:A:H2'	36:1:2697:A:C8	2.55	0.42
36:1:2374:C:N4	36:1:2941:A:N3	2.67	0.42
36:1:3001:C:H2'	36:1:3002:C:H6	1.85	0.42
36:1:3380:U:H2'	36:1:3381:U:C6	2.54	0.42
1:2:1280:C:H1'	1:2:1429:G:N2	2.35	0.42
1:2:397:A:H5''	10:S8:51:GLY:H	1.85	0.42
1:2:373:G:N2	1:2:604:A:OP1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:77:U:H4'	1:2:78:A:O5'	2.19	0.42
1:2:848:C:H2'	1:2:849:C:H6	1.85	0.42
1:2:947:U:OP1	3:S1:165:ARG:NE	2.44	0.42
37:3:64:A:N6	47:M0:200:LEU:HD11	2.34	0.42
36:5:1461:A:H2'	36:5:1462:A:O4'	2.20	0.42
36:5:1562:C:H2'	36:5:1563:C:H6	1.85	0.42
36:5:2255:A:HO2'	36:5:2256:A:P	2.43	0.42
36:5:2255:A:O2'	36:5:2256:A:OP 2	2.35	0.42
36:5:2409:G:H4'	36:5:2410:U:OP 2	2.20	0.42
36:5:254:A:H2'	36:5:255:A:C8	2.55	0.42
68:O2:50:ILE:HG22	36:5:426:G:H5'	192.27	0.42
36:5:787:G:H2'	36:5:788:C:C6	2.55	0.42
1:6:1227:A:H4'	1:6:1228:G:H5'	2.01	0.42
1:6:1299:G:H2'	1:6:1300:A:C8	2.54	0.42
1:6:1297:G:N2	1:6:1300:A:OP2	2.48	0.42
1:6:1325:A:H2'	1:6:1326:A:C8	2.54	0.42
32:E0:28:LYS:HZ1	1:6:542:A:H61	428.21	0.42
1:6:775:G:N2	1:6:786:C:N3	2.68	0.42
1:6:839:U:H2'	1:6:840:U:H6	1.84	0.42
1:6:947:U:H2'	1:6:948:G:H8	1.84	0.42
1:6:97:C:H2'	1:6:98:U:H6	1.85	0.42
12:C0:38:LYS:HD3	12:C0:41:TYR:OH	2.19	0.42
13:C1:22:ASN:ND2	13:C1:25:VAL:HG23	2.34	0.42
1:2:1229:G:O6	14:C2:47:GLU:HG3	2.20	0.42
15:C3:56:ASP:OD2	29:D7:52:THR:OG1	4.22	0.42
16:C4:24:ASN:N	16:C4:55:SER:HB3	2.35	0.42
18:C6:8:GLN:HA	18:C6:21:HIS:HA	2.02	0.42
20:C8:127:HIS:CD2	20:C8:133:VAL:HG21	2.54	0.42
20:C8:49:LYS:NZ	20:C8:80:LYS:HB2	6.06	0.42
23:D1:12:TYR:O	23:D1:14:PRO:HD3	2.19	0.42
2:S0:63:ILE:HG23	23:D1:35:ASN:O	2.19	0.42
23:D1:71:ARG:HB2	23:D1:83:TRP:NE1	4.89	0.42
24:D2:5:SER:C	24:D2:7:LEU:N	3.32	0.42
39:L2:206:PRO:HG3	39:L2:213:GLY:CA	2.72	0.42
39:L2:44:ILE:O	39:L2:61:VAL:HG23	4.87	0.42
40:L3:215:ILE:HG12	40:L3:280:HIS:O	2.19	0.42
40:L3:232:ARG:HD2	40:L3:269:GLN:O	2.76	0.42
41:L4:33:ASP:O	41:L4:37:THR:HG23	2.19	0.42
44:L7:130:ILE:O	44:L7:134:VAL:HG22	2.20	0.42
38:4:155:A:H5'	45:L8:185:ARG:NH2	2.35	0.42
45:L8:24:ASN:N	45:L8:25:PRO:HD2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:M0:98:ARG:HB3	47:M0:120:GLY:HA3	2.02	0.42
50:M4:43:LYS:NZ	50:M4:43:LYS:HB2	2.34	0.42
55:M9:175:GLN:HG3	55:M9:176:ARG:HD3	4.61	0.42
56:N0:115:ARG:N	56:N0:115:ARG:HD2	2.35	0.42
56:N0:155:ARG:NH2	56:N0:172:TYR:H	5.76	0.42
62:N6:39:LEU:HA	62:N6:39:LEU:HD23	1.76	0.42
62:N6:54:ASP:HB2	62:N6:70:ILE:HD13	3.11	0.42
62:N6:80:VAL:HG12	62:N6:101:PRO:HD3	2.01	0.42
63:N7:46:ILE:HD11	63:N7:48:ARG:C	2.40	0.42
63:N7:86:THR:O	63:N7:127:ASN:ND2	4.92	0.42
64:N8:120:ASN:OD1	64:N8:141:ALA:HA	3.04	0.42
64:N8:21:ARG:HD2	36:5:1369:A:C5'	185.39	0.42
64:N8:24:LYS:HD2	64:N8:26:ARG:NE	2.33	0.42
49:M3:64:LYS:HD2	64:N8:66:ALA:HB1	3.28	0.42
72:O6:52:PRO:HA	72:O6:55:ARG:NH1	2.35	0.42
72:O6:66:GLU:O	72:O6:70:ARG:HG3	2.20	0.42
74:O8:64:LYS:HG3	74:O8:65:LEU:N	4.07	0.42
2:S0:56:LYS:HD2	2:S0:158:VAL:HG23	2.01	0.42
4:S2:122:ALA:O	4:S2:126:ARG:HG3	2.20	0.42
5:S3:37:VAL:HG12	5:S3:50:ILE:HA	2.44	0.42
5:S3:42:THR:OG1	5:S3:44:THR:O	5.95	0.42
6:S4:122:LYS:HE3	6:S4:122:LYS:HB2	2.46	0.42
6:S4:176:ASP:HB2	6:S4:179:LYS:HE2	2.02	0.42
6:S4:49:ARG:HD2	6:S4:56:LEU:O	2.19	0.42
7:S5:109:LYS:O	7:S5:113:ILE:HG13	2.31	0.42
7:S5:158:GLN:HG2	30:D8:66:LEU:HD11	2.02	0.42
9:S7:63:PRO:O	9:S7:64:VAL:HB	2.20	0.42
11:S9:66:ASP:OD2	11:S9:68:LYS:N	2.79	0.42
34:SR:81:LEU:HD22	34:SR:113:VAL:HG21	2.00	0.42
36:1:1176:C:OP1	52:M6:25:LYS:HE3	2.19	0.42
36:1:1599:G:C6	36:1:1600:U:C4	3.08	0.42
36:1:256:G:H2'	36:1:257:U:C6	2.55	0.42
36:1:1132:C:H4'	36:1:2865:U:O2'	2.19	0.42
36:1:3267:A:H2'	43:L6:69:PHE:CE1	2.55	0.42
36:1:3289:G:C2	36:1:3290:G:C8	3.08	0.42
36:1:3333:G:N2	36:1:3369:G:O2'	2.53	0.42
36:1:511:G:H2'	36:1:512:U:O4'	2.19	0.42
1:2:1120:U:H2'	1:2:1121:C:C6	2.55	0.42
1:2:1186:U:OP2	1:2:1456:C:H1'	2.19	0.42
1:2:1435:G:O6	12:C0:64:TYR:OH	2.31	0.42
1:2:1654:G:C6	1:2:1745:G:C6	3.08	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:434:G:N2	1:2:436:A:H3'	2.35	0.42
1:2:539:G:OP2	1:2:539:G:H8	2.03	0.42
37:3:7:G:H5''	42:L5:22:ARG:HD3	2.02	0.42
38:4:121:U:C2	38:4:122:U:C5	3.08	0.42
36:5:119:U:H4'	36:5:120:G:H3'	2.01	0.42
36:5:1252:A:H2'	36:5:1253:U:H5'	2.02	0.42
36:5:1303:A:O4'	36:5:2885:C:O2'	2.38	0.42
36:5:1113:G:O2'	36:5:1369:A:N3	2.47	0.42
36:5:2309:A:OP1	36:5:2309:A:H8	2.03	0.42
36:5:2658:G:C6	36:5:2659:G:N7	2.88	0.42
36:5:2765:C:H2'	36:5:2766:U:C6	2.54	0.42
36:5:2946:A:H5''	36:5:2947:G:H5'	2.02	0.42
36:5:3018:C:C4	36:5:3019:U:C4	3.07	0.42
36:5:1940:G:N2	36:5:3362:A:H8	2.13	0.42
36:5:767:U:O2'	36:5:768:C:H5''	2.20	0.42
1:6:1153:G:H2'	1:6:1154:G:O4'	2.19	0.42
1:6:1358:G:H1	1:6:1365:C:H42	1.68	0.42
8:S6:176:GLN:HG2	1:6:169:A:C5'	327.62	0.42
1:6:1756[A]:A:OP2	1:6:1756[A]:A:H8	2.03	0.42
1:6:526:A:N6	1:6:527:A:C6	2.88	0.42
1:6:583:C:H2'	1:6:584:C:H6	1.84	0.42
6:S4:21:ASP:HB2	1:6:773:C:OP1	388.10	0.42
1:6:836:U:H2'	1:6:837:G:C8	2.55	0.42
37:7:94:C:H2'	37:7:95:A:C8	2.55	0.42
38:8:79:A:H2'	38:8:80:A:O4'	2.19	0.42
13:C1:74:THR:HG23	13:C1:122:ILE:HG13	2.02	0.42
15:C3:42:ARG:O	15:C3:44:GLY:N	3.23	0.42
15:C3:74:ILE:H	15:C3:74:ILE:HG13	3.60	0.42
16:C4:20:TYR:HA	16:C4:84:ARG:HG2	4.11	0.42
17:C5:127:ARG:O	17:C5:129:GLY:N	4.41	0.42
20:C8:29:VAL:HB	20:C8:30:TYR:CD1	2.96	0.42
20:C8:82:PRO:HD2	20:C8:85:PHE:HB2	2.02	0.42
21:C9:16:ASN:O	21:C9:19:ALA:HB3	2.20	0.42
26:D4:7:ILE:HD11	26:D4:43:LYS:HD2	2.01	0.42
28:D6:53:LEU:HD13	28:D6:53:LEU:HA	3.11	0.42
7:S5:144:GLU:CD	30:D8:57:MET:HG3	4.67	0.42
1:2:1253:U:H4'	33:E1:143:LYS:N	2.35	0.42
39:L2:11:GLY:HA3	36:5:2163:C:O2'	179.16	0.42
39:L2:188:LYS:O	39:L2:192:LYS:HD2	2.19	0.42
39:L2:195:SER:O	39:L2:198:LYS:NZ	2.84	0.42
39:L2:32:LEU:HD23	39:L2:163:ARG:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:L4:55:LYS:HG2	41:L4:59:GLN:NE2	4.50	0.42
42:L5:258:LYS:HD2	42:L5:265:TYR:HE2	1.85	0.42
43:L6:90:LYS:HE2	43:L6:90:LYS:HB3	1.77	0.42
45:L8:68:ARG:HG2	45:L8:68:ARG:H	2.16	0.42
46:L9:174:LYS:HB2	76:Q0:127:LEU:HD11	2.02	0.42
47:M0:30:LYS:HD3	47:M0:30:LYS:HA	4.21	0.42
49:M3:179:PHE:HB2	49:M3:183:ARG:NH1	3.07	0.42
50:M4:122:VAL:O	50:M4:126:GLN:HG3	2.36	0.42
51:M5:55:ALA:HB3	36:5:149:U:H5'	100.38	0.42
52:M6:129:LEU:HD12	52:M6:129:LEU:HA	1.78	0.42
53:M7:136:ILE:HD11	36:5:1846:C:C2	141.41	0.42
54:M8:170:ARG:O	54:M8:171:LYS:HG2	2.20	0.42
41:L4:30:ILE:N	54:M8:25:TYR:OH	2.53	0.42
55:M9:109:TYR:HB3	55:M9:115:ILE:HG22	2.01	0.42
55:M9:88:ARG:HB3	55:M9:88:ARG:HH11	4.06	0.42
36:1:1722:U:H1'	55:M9:96:ILE:HG12	2.02	0.42
55:M9:9:ARG:NH2	55:M9:10:LEU:HD13	2.35	0.42
58:N2:80:THR:HG21	58:N2:95:PHE:HD2	6.36	0.42
59:N3:34:LEU:HA	59:N3:34:LEU:HD23	3.34	0.42
62:N6:118:LEU:HD13	62:N6:121:ARG:HH11	3.21	0.42
63:N7:4:PHE:CE2	66:O0:35:ARG:HA	2.55	0.42
63:N7:51:LEU:HD12	63:N7:65:ARG:HD2	2.01	0.42
64:N8:7:LYS:HA	64:N8:7:LYS:HD3	1.89	0.42
65:N9:59:LYS:HE3	65:N9:59:LYS:HB2	4.66	0.42
67:O1:74:ARG:NH1	67:O1:109:VAL:HG11	5.00	0.42
68:O2:79:VAL:HG13	68:O2:111:ARG:HG2	2.01	0.42
70:O4:44:CYS:HG	70:O4:47:CYS:HG	2.15	0.42
71:O5:85:THR:HB	71:O5:88:LEU:HB2	2.00	0.42
72:O6:43:LEU:O	72:O6:43:LEU:HD22	2.66	0.42
74:O8:70:PRO:HA	74:O8:71:PRO:HD3	1.82	0.42
79:Q3:79:VAL:O	79:Q3:83:ILE:HG12	2.52	0.42
2:S0:142:PRO:HB3	23:D1:34:ILE:CD1	2.50	0.42
2:S0:41:ARG:NH1	2:S0:45:VAL:HG23	2.35	0.42
2:S0:54:TRP:O	2:S0:58:VAL:HG23	2.20	0.42
3:S1:83:LYS:HE3	3:S1:83:LYS:HB2	1.77	0.42
5:S3:173:ARG:HD3	5:S3:173:ARG:HA	1.92	0.42
5:S3:210:GLU:HA	5:S3:211:PRO:HD3	1.91	0.42
6:S4:166:SER:O	6:S4:168:LYS:HG2	5.40	0.42
6:S4:206:ASP:N	6:S4:206:ASP:OD1	2.53	0.42
7:S5:112:ARG:HD2	18:C6:43:ILE:HD11	2.02	0.42
7:S5:200:ASN:HB3	7:S5:208:SER:HB2	4.09	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:173:TYR:CD1	9:S7:181:ILE:HD11	4.29	0.42
11:S9:128:LEU:O	11:S9:133:HIS:HB2	2.20	0.42
11:S9:156:ILE:HG13	11:S9:156:ILE:H	1.77	0.42
11:S9:54:ARG:HA	11:S9:57:ARG:HE	1.93	0.42
34:SR:248:ASN:ND2	34:SR:298:GLY:HA3	3.40	0.42
34:SR:61:PHE:HB3	34:SR:92:TRP:CZ3	2.92	0.42
36:1:104:G:H2'	36:1:105:C:O4'	2.20	0.42
36:1:1274:A:H2'	36:1:1275:C:C6	2.54	0.42
36:1:1394:A:N3	38:4:19:C:O2'	2.42	0.42
36:1:2202:C:H2'	36:1:2203:U:O4'	2.20	0.42
36:1:2561:A:O2'	36:1:2562:A:O5'	2.33	0.42
36:1:3138:U:OP2	40:L3:30:LYS:HD3	2.20	0.42
36:1:66:A:O2'	36:1:315:C:O2'	2.32	0.42
36:1:394:G:N2	36:1:396:A:H3'	2.35	0.42
36:1:585:A:H5''	69:O3:70:LYS:CE	2.49	0.42
1:2:1147:A:O2'	1:2:1635:A:H2'	2.20	0.42
1:2:1325:A:C2	1:2:1326:A:C5	3.08	0.42
1:2:1524:A:H2	1:2:1590:G:H1'	1.85	0.42
1:2:40:A:H2'	1:2:41:A:O4'	2.20	0.42
1:2:542:A:N7	1:2:543:C:H2'	2.35	0.42
37:3:112:G:H2'	37:3:113:C:H6	1.85	0.42
36:5:1027:A:H4'	36:5:1028:U:OP2	2.18	0.42
36:5:1087:G:C2	36:5:1088:U:C4	3.08	0.42
36:5:1108:U:H2'	36:5:1109:U:H6	1.83	0.42
36:5:1340:G:C6	36:5:1341:U:C4	3.08	0.42
36:5:209:A:H4'	36:5:211:A:H8	1.82	0.42
36:5:2448:G:N2	36:5:2499:U:H3	2.18	0.42
54:M8:107:THR:HG21	36:5:676:G:H3'	136.44	0.42
36:5:747:A:C6	36:5:748:U:C4	3.08	0.42
36:5:850:U:H2'	36:5:851:C:H6	1.85	0.42
1:6:116:U:O2	1:6:333:A:H2	2.03	0.42
1:6:1195:C:H5''	1:6:1197:C:C6	2.55	0.42
1:6:1516:A:O2'	1:6:1517:U:H5'	2.20	0.42
1:6:329:G:H2'	1:6:330:G:C8	2.55	0.42
1:6:649:U:H2'	1:6:650:U:H5	1.85	0.42
38:8:40:A:H2'	38:8:41:A:H8	1.85	0.42
38:8:6:U:H2'	38:8:7:U:C6	2.54	0.42
13:C1:57:LYS:O	13:C1:138:ASN:ND2	2.50	0.42
17:C5:24:LYS:H	17:C5:24:LYS:HD2	1.85	0.42
18:C6:86:ALA:HB1	18:C6:109:PHE:CE2	2.77	0.42
19:C7:60:ARG:HG3	19:C7:66:VAL:HG21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:C9:141:GLU:O	21:C9:143:ASP:N	3.56	0.42
39:L2:44:ILE:HG23	39:L2:87:PHE:CE1	3.74	0.42
36:1:1651:U:C5'	39:L2:71:LEU:HD13	2.50	0.42
40:L3:240:ARG:HA	40:L3:246:LEU:HD22	5.51	0.42
36:1:2947:G:N3	40:L3:250:ALA:HB1	2.35	0.42
42:L5:115:LEU:HD12	42:L5:119:TYR:CD2	4.59	0.42
42:L5:218:ARG:HA	42:L5:221:GLU:OE2	2.20	0.42
42:L5:41:LYS:HD3	42:L5:41:LYS:HA	3.84	0.42
44:L7:163:LEU:HD21	44:L7:201:PHE:CE1	2.55	0.42
44:L7:214:TRP:CE2	44:L7:219:LYS:HE3	2.55	0.42
44:L7:95:ILE:HA	44:L7:96:PRO:HD3	1.92	0.42
45:L8:128:LYS:HA	45:L8:129:PRO:HD3	1.82	0.42
45:L8:98:ARG:HE	45:L8:189:LEU:HA	1.85	0.42
46:L9:49:ASN:HD21	46:L9:52:LEU:HB2	1.85	0.42
47:M0:86:HIS:HB3	47:M0:139:ARG:HG2	2.02	0.42
47:M0:65:LEU:HD23	47:M0:159:PHE:CE1	2.55	0.42
48:M1:37:LEU:HA	48:M1:37:LEU:HD23	2.34	0.42
50:M4:20:VAL:HA	50:M4:33:ALA:O	2.20	0.42
52:M6:121:PRO:HA	52:M6:124:LEU:HD22	2.63	0.42
53:M7:137:ASN:HD22	53:M7:137:ASN:HA	1.56	0.42
53:M7:139:TYR:CD2	36:5:2355:G:H4'	146.94	0.42
36:1:412:G:H5'	53:M7:26:PHE:HZ	1.84	0.42
57:N1:18:ASP:HB2	57:N1:21:LYS:HB2	2.78	0.42
58:N2:56:VAL:HG22	58:N2:65:VAL:HG13	2.02	0.42
59:N3:39:VAL:HG22	59:N3:52:ALA:HB2	2.02	0.42
59:N3:74:MET:SD	59:N3:102:ILE:HD13	2.60	0.42
64:N8:16:SER:HA	36:5:942:U:H3	170.53	0.42
76:Q0:104:PRO:HA	76:Q0:105:PRO:HD3	1.79	0.42
2:S0:85:ALA:HA	2:S0:202:TYR:HD1	1.84	0.42
3:S1:36:SER:HB3	3:S1:41:ARG:NH2	8.39	0.42
3:S1:81:PHE:HB3	3:S1:106:THR:OG1	2.20	0.42
4:S2:218:ILE:H	4:S2:218:ILE:HG13	1.70	0.42
4:S2:77:GLN:N	4:S2:77:GLN:HE21	2.15	0.42
4:S2:89:GLN:HG3	4:S2:94:GLN:HG2	2.01	0.42
7:S5:42:LEU:HA	7:S5:48:PHE:HD2	1.85	0.42
8:S6:13:GLN:HE21	8:S6:13:GLN:HB3	1.55	0.42
9:S7:102:PRO:HD3	9:S7:112:ARG:HD3	2.45	0.42
9:S7:150:GLN:HB3	9:S7:181:ILE:HD12	2.02	0.42
9:S7:46:ILE:HG12	9:S7:60:ILE:HA	2.20	0.42
11:S9:114:TYR:O	11:S9:116:LEU:N	3.02	0.42
35:SM:50:ASN:HB2	35:SM:51:ARG:H	4.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:SR:248:ASN:ND2	34:SR:249:ARG:HG3	2.35	0.42
34:SR:255:ALA:HB2	34:SR:292:LEU:HD22	2.02	0.42
36:1:1673:G:C4	36:1:1775:G:C2	3.08	0.41
36:1:40:A:N7	64:N8:29:PRO:O	2.53	0.41
36:1:437:G:N2	36:1:622:A:H61	2.18	0.41
36:1:681:U:O4	41:L4:118:LYS:NZ	2.38	0.41
1:2:621:A:N3	1:2:1107:G:H1'	2.35	0.41
1:2:144:U:O2'	1:2:145:A:H8	2.02	0.41
1:2:1486:G:C6	1:2:1522:U:C5	3.07	0.41
1:2:1575:G:H2'	1:2:1576:A:C8	2.55	0.41
1:2:1615:C:C4	7:S5:81:ARG:HA	2.55	0.41
1:2:979:A:N3	1:2:1775:U:O2'	2.53	0.41
37:3:37:G:C2	37:3:38:U:C2	3.08	0.41
38:4:97:A:OP1	71:O5:67:ARG:NH2	2.38	0.41
36:5:1240:A:N3	36:5:1249:G:N2	2.68	0.41
36:5:1549:U:H2'	36:5:1550:C:H6	1.85	0.41
36:5:2266:U:O2'	36:5:2267:C:H5'	2.20	0.41
49:M3:45:LYS:HE2	36:5:241:G:O4'	45.57	0.41
36:5:2772:C:H1'	36:5:2773:C:OP2	2.20	0.41
36:5:1205:A:H4'	36:5:2835:U:O2'	2.19	0.41
36:5:2849:C:H2'	36:5:2850:G:O4'	2.20	0.41
40:L3:128:LYS:HG3	36:5:3294:A:H5'	197.66	0.41
8:S6:96:SER:OG	1:6:420:A:OP1	295.84	0.41
42:L5:269:SER:OG	37:7:1:G:N3	315.78	0.41
13:C1:109:VAL:HG23	13:C1:137:PHE:C	2.40	0.41
13:C1:26:LYS:HD3	13:C1:26:LYS:HA	1.80	0.41
16:C4:56:SER:HA	16:C4:57:PRO:HD3	1.89	0.41
19:C7:49:LYS:HA	1:6:1389:C:H4'	421.96	0.41
19:C7:71:PHE:O	19:C7:73:LEU:N	2.50	0.41
1:2:1477:G:H5''	21:C9:45:MET:O	2.19	0.41
21:C9:58:ALA:HB1	21:C9:108:LEU:HD11	2.02	0.41
22:D0:72:ASN:HA	1:6:1198:G:O2'	386.62	0.41
4:S2:222:TYR:OH	23:D1:11:LEU:O	2.36	0.41
24:D2:32:LYS:HE3	24:D2:32:LYS:HB2	1.88	0.41
25:D3:92:CYS:O	25:D3:95:PHE:HB2	3.01	0.41
26:D4:126:ALA:HA	26:D4:129:VAL:HG12	2.03	0.41
39:L2:66:PRO:HB2	39:L2:67:TYR:CE2	2.68	0.41
40:L3:15:GLY:O	36:5:3009:G:N2	241.35	0.41
42:L5:173:VAL:HA	42:L5:174:PRO:HD2	2.01	0.41
42:L5:54:ARG:HG3	37:7:5:G:O3'	281.95	0.41
41:L4:311:HIS:CE1	44:L7:162:PRO:HG2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:L9:154:VAL:O	46:L9:158:ALA:N	2.42	0.41
46:L9:18:VAL:HG12	46:L9:27:VAL:HG22	2.03	0.41
48:M1:105:GLY:HA3	36:5:2674:A:H5''	332.76	0.41
49:M3:168:ARG:O	49:M3:168:ARG:HG3	3.38	0.41
50:M4:135:LEU:O	50:M4:137:LYS:N	2.82	0.41
51:M5:172:ARG:HD2	36:5:30:G:O5'	111.40	0.41
51:M5:199:LEU:HB3	51:M5:203:ARG:HE	1.85	0.41
51:M5:19:LEU:HA	51:M5:19:LEU:HD12	1.78	0.41
51:M5:35:VAL:O	51:M5:64:VAL:HA	2.19	0.41
36:1:115:A:P	51:M5:49:ARG:HE	2.43	0.41
52:M6:41:LEU:HB3	52:M6:138:LEU:HD22	2.01	0.41
53:M7:3:ARG:O	53:M7:18:ARG:NH2	4.12	0.41
56:N0:1:MET:SD	56:N0:36:ILE:HG21	2.59	0.41
57:N1:116:ARG:HA	57:N1:126:VAL:HG11	3.39	0.41
58:N2:33:TYR:CE2	58:N2:63:VAL:HG21	2.55	0.41
58:N2:76:LEU:HA	58:N2:79:LEU:HD12	2.01	0.41
59:N3:135:VAL:HG11	60:N4:26:SER:HB3	2.01	0.41
61:N5:106:ASP:OD2	61:N5:107:VAL:N	5.09	0.41
61:N5:65:GLN:HB2	61:N5:65:GLN:HE21	3.98	0.41
62:N6:57:LEU:HB3	62:N6:105:VAL:CG1	3.48	0.41
64:N8:103:ASP:HA	64:N8:126:LYS:HB2	2.01	0.41
66:O0:26:GLY:O	66:O0:89:VAL:HG13	2.20	0.41
70:O4:11:ASN:HA	70:O4:12:PRO:HD3	1.74	0.41
73:O7:68:LYS:HD2	73:O7:69:HIS:NE2	2.35	0.41
7:S5:141:GLY:HA2	7:S5:142:PRO:HD3	1.91	0.41
7:S5:28:PRO:HB2	7:S5:29:ILE:H	3.36	0.41
8:S6:173:PRO:HG3	1:6:66:U:C6	334.61	0.41
10:S8:48:THR:HG21	10:S8:54:LYS:HG3	2.01	0.41
36:1:1326:A:H2'	36:1:1327:C:O4'	2.20	0.41
36:1:2917:G:H4'	59:N3:48:ARG:O	2.19	0.41
36:1:594:U:H2'	36:1:609:G:O6	2.19	0.41
36:1:674:G:O6	54:M8:56:LYS:NZ	2.50	0.41
36:1:66:A:N6	36:1:75:G:N2	2.69	0.41
1:2:1486:G:C8	1:2:1487:A:C8	3.08	0.41
1:2:766:U:H3'	1:2:768:C:OP2	2.20	0.41
1:2:866:G:N2	1:2:965:U:C5	2.88	0.41
37:3:25:G:H2'	37:3:26:C:O4'	2.20	0.41
38:4:86:U:H3'	71:O5:7:TYR:OH	2.19	0.41
46:L9:62:ARG:HD2	36:5:1211:U:OP2	319.99	0.41
36:5:915:A:H8	36:5:2136:C:O2'	2.03	0.41
36:5:2542:U:O2'	36:5:2543:U:O5'	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:3351:U:H5'	36:5:3352:U:OP2	2.20	0.41
36:5:370:U:H4'	36:5:404:G:H5'	2.02	0.41
36:5:411:U:H2'	36:5:412:G:C8	2.54	0.41
1:6:1119:G:H2'	1:6:1120:U:O4'	2.20	0.41
14:C2:45:LEU:HB2	1:6:1228:G:OP1	462.70	0.41
1:6:1241:G:H2'	1:6:1242:A:O4'	2.20	0.41
1:6:1437:U:H2'	1:6:1438:G:C8	2.55	0.41
1:6:1640:C:H1'	1:6:1763:A:N1	2.34	0.41
1:6:436:A:H8	1:6:436:A:O5'	2.03	0.41
12:C0:24:LYS:HB2	12:C0:63:TYR:CE1	2.54	0.41
12:C0:50:THR:HG22	12:C0:55:VAL:HG13	6.54	0.41
15:C3:34:ILE:HG13	15:C3:67:THR:HG21	2.00	0.41
15:C3:46:THR:HG21	15:C3:90:TYR:OH	3.91	0.41
18:C6:102:LYS:HE2	18:C6:102:LYS:HB3	1.83	0.41
20:C8:100:THR:HG21	20:C8:108:LYS:HG2	2.03	0.41
20:C8:49:LYS:HG3	20:C8:81:ILE:HD11	3.77	0.41
22:D0:97:VAL:HG13	22:D0:98:GLN:H	2.17	0.41
4:S2:140:ARG:HA	23:D1:10:GLU:OE1	2.20	0.41
24:D2:24:GLN:OE1	29:D7:4:VAL:HA	2.20	0.41
25:D3:103:LEU:HD12	25:D3:126:LYS:HE2	2.01	0.41
25:D3:57:LEU:HD23	25:D3:57:LEU:HA	1.73	0.41
25:D3:69:ARG:HD3	25:D3:117:ILE:CG2	2.51	0.41
26:D4:37:LYS:HD2	26:D4:57:VAL:HG23	6.74	0.41
29:D7:16:ALA:HA	29:D7:29:ARG:HH21	4.22	0.41
39:L2:133:TYR:HB3	39:L2:168:VAL:HB	2.97	0.41
39:L2:70:ARG:NH2	39:L2:72:ARG:HH21	6.19	0.41
40:L3:216:ASP:CG	40:L3:278:ILE:HG22	2.41	0.41
41:L4:118:LYS:O	41:L4:122:THR:HG23	2.69	0.41
42:L5:51:LEU:HB2	42:L5:144:VAL:HG13	3.00	0.41
42:L5:211:LEU:HD23	42:L5:211:LEU:HA	2.57	0.41
42:L5:205:SER:HB3	42:L5:236:LEU:HD22	2.67	0.41
42:L5:294:ALA:C	42:L5:296:GLN:H	2.23	0.41
44:L7:98:LYS:HE2	44:L7:129:LEU:HD21	2.01	0.41
45:L8:78:PHE:O	45:L8:80:TYR:N	2.43	0.41
46:L9:17:THR:HG23	46:L9:28:VAL:HB	5.05	0.41
47:M0:95:HIS:HB2	47:M0:128:ARG:HD2	2.94	0.41
47:M0:159:PHE:HA	47:M0:160:PRO:HD2	2.50	0.41
47:M0:188:GLY:HA3	47:M0:216:TYR:HD1	1.85	0.41
49:M3:28:GLN:HB3	51:M5:201:ARG:NH1	2.99	0.41
51:M5:28:TRP:O	51:M5:32:GLN:HG2	2.21	0.41
52:M6:48:PHE:O	52:M6:52:LEU:HG	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:M7:126:ARG:HA	53:M7:140:GLU:HG2	2.39	0.41
54:M8:115:VAL:O	54:M8:118:GLY:N	2.89	0.41
54:M8:161:LYS:O	54:M8:162:ALA:HB3	2.19	0.41
36:1:522:A:O2'	56:N0:66:GLU:O	2.37	0.41
57:N1:66:ASN:OD1	57:N1:67:VAL:N	3.14	0.41
58:N2:37:LEU:O	58:N2:41:ILE:HG13	2.20	0.41
60:N4:39:LEU:HA	60:N4:39:LEU:HD12	1.82	0.41
57:N1:84:TYR:CD1	65:N9:23:LYS:HG3	6.92	0.41
66:O0:27:TYR:O	66:O0:31:VAL:HG23	2.20	0.41
69:O3:37:THR:OG1	69:O3:39:GLN:OE1	2.37	0.41
69:O3:89:LEU:HA	69:O3:89:LEU:HD23	1.87	0.41
69:O3:8:TYR:HB3	69:O3:101:PHE:CE1	2.88	0.41
72:O6:57:LEU:HD22	72:O6:57:LEU:HA	2.88	0.41
73:O7:21:ARG:HD2	73:O7:37:CYS:SG	2.78	0.41
73:O7:19:CYS:SG	73:O7:34:CYS:CB	3.03	0.41
2:S0:152:PRO:C	2:S0:154:GLU:H	2.22	0.41
2:S0:56:LYS:O	2:S0:60:ALA:N	2.85	0.41
2:S0:71:GLU:HA	2:S0:94:GLY:C	3.01	0.41
3:S1:103:MET:HB3	3:S1:215:VAL:CG2	4.66	0.41
4:S2:107:SER:O	4:S2:192:GLY:HA2	2.20	0.41
4:S2:53:ILE:HG23	4:S2:56:ILE:HD11	2.01	0.41
4:S2:84:LYS:HA	4:S2:85:PRO:HD3	1.83	0.41
4:S2:88:LYS:HE2	4:S2:90:THR:HB	2.01	0.41
5:S3:21:LEU:HD23	5:S3:21:LEU:HA	1.87	0.41
6:S4:206:ASP:O	6:S4:222:LEU:N	2.47	0.41
7:S5:58:LEU:HD13	7:S5:138:THR:HG22	2.03	0.41
36:1:2310:U:H6	36:1:2310:U:O5'	2.03	0.41
36:1:2623:G:C4	36:1:2624:G:C8	3.09	0.41
36:1:2735:U:H2'	36:1:2736:A:C8	2.56	0.41
36:1:290:G:H5''	51:M5:98:LEU:HD23	2.02	0.41
36:1:312:C:H2'	36:1:313:A:C8	2.55	0.41
36:1:439:C:H3'	36:1:440:A:H8	1.85	0.41
1:2:1166:A:H2'	1:2:1167:G:O4'	2.21	0.41
1:2:1727:G:H2'	1:2:1728:A:C8	2.55	0.41
1:2:736:C:OP1	6:S4:197:HIS:NE2	2.52	0.41
1:2:738:G:H2'	1:2:739:G:C8	2.48	0.41
36:5:1128:U:H2'	36:5:1129:A:O4'	2.21	0.41
36:5:1200:A:H5'	36:5:1201:C:OP1	2.21	0.41
36:5:1222:G:OP2	36:5:1222:G:H8	2.03	0.41
36:5:1464:G:N1	36:5:1467:A:OP2	2.53	0.41
36:5:1733:G:H2'	36:5:1734:G:H8	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:174:C:H2'	36:5:175:C:C6	2.54	0.41
36:5:1782:U:H2'	36:5:1783:U:O4'	2.21	0.41
36:5:2097:U:H2'	36:5:2098:C:H6	1.85	0.41
36:5:2514:U:H6	36:5:2514:U:OP1	2.03	0.41
57:N1:6:GLY:HA2	36:5:2630:C:OP1	230.33	0.41
36:5:2980:U:H2'	36:5:2981:U:H6	1.84	0.41
36:5:3060:C:H2'	36:5:3061:G:C8	2.55	0.41
1:6:1239:U:H2'	1:6:1240:U:C6	2.54	0.41
1:6:1778:G:H2'	1:6:1779:U:H6	1.85	0.41
1:6:533:U:H2'	1:6:534:A:H5''	2.01	0.41
1:6:994:G:H2'	1:6:995:A:H8	1.85	0.41
38:8:72:A:C2	38:8:89:A:H5'	2.55	0.41
18:C6:60:PHE:CE2	18:C6:89:LEU:HD22	4.44	0.41
19:C7:5:ARG:HB2	19:C7:10:LYS:HE2	2.03	0.41
19:C7:17:ILE:HD11	19:C7:54:THR:O	2.20	0.41
19:C7:87:GLU:O	19:C7:88:VAL:HG12	2.21	0.41
20:C8:78:HIS:O	20:C8:80:LYS:N	3.26	0.41
21:C9:126:GLU:HG2	21:C9:126:GLU:H	3.29	0.41
21:C9:57:ARG:NH2	21:C9:80:TYR:HB3	2.35	0.41
23:D1:44:ARG:HE	23:D1:44:ARG:HB2	1.72	0.41
25:D3:13:ARG:HA	25:D3:16:ARG:NH2	3.91	0.41
39:L2:174:ARG:HH22	36:5:2180:G:P	212.22	0.41
40:L3:287:LYS:HA	40:L3:320:ASP:OD1	2.98	0.41
40:L3:35:ASP:HA	40:L3:184:ASN:ND2	3.05	0.41
40:L3:60:LEU:HD11	40:L3:62:ARG:HB2	2.61	0.41
41:L4:185:LYS:HE2	41:L4:199:TRP:HB3	2.76	0.41
41:L4:219:LEU:HD23	41:L4:222:VAL:HG11	2.66	0.41
41:L4:317:PRO:HB2	44:L7:149:TYR:HD1	1.84	0.41
41:L4:67:THR:OG1	36:5:2402:A:H5''	173.22	0.41
41:L4:99:MET:HE2	41:L4:103:THR:H	3.93	0.41
42:L5:233:ALA:O	42:L5:236:LEU:HB2	2.28	0.41
43:L6:145:LEU:O	43:L6:148:GLU:N	2.52	0.41
46:L9:43:VAL:CG2	46:L9:55:VAL:HG12	3.25	0.41
49:M3:188:ARG:NH2	49:M3:189:GLU:HG3	2.35	0.41
49:M3:54:LEU:HD13	49:M3:75:PHE:CZ	2.56	0.41
49:M3:95:ILE:HD13	49:M3:116:LEU:HD22	2.03	0.41
50:M4:113:THR:N	50:M4:116:GLU:HB3	4.14	0.41
51:M5:183:THR:O	51:M5:183:THR:OG1	2.90	0.41
54:M8:8:LYS:HB2	54:M8:8:LYS:HE3	1.65	0.41
55:M9:155:LEU:HA	55:M9:158:GLU:HG2	2.03	0.41
55:M9:180:LYS:HE3	55:M9:180:LYS:HB3	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:3333:G:C8	60:N4:51:TRP:CD1	3.08	0.41
61:N5:91:ASN:O	61:N5:94:GLN:N	3.08	0.41
63:N7:129:TRP:O	63:N7:131:PHE:N	3.45	0.41
64:N8:79:TRP:CZ2	64:N8:123:VAL:HG13	5.22	0.41
68:O2:119:VAL:O	68:O2:122:PRO:HD3	2.41	0.41
68:O2:20:HIS:CG	68:O2:42:VAL:HG21	2.55	0.41
70:O4:59:PRO:HD3	36:5:1654:A:O2'	167.58	0.41
74:O8:46:ARG:NH1	74:O8:47:GLY:O	2.54	0.41
2:S0:41:ARG:HE	2:S0:45:VAL:HG21	5.12	0.41
3:S1:120:LEU:HD22	3:S1:122:GLU:HG3	2.71	0.41
3:S1:158:SER:HA	3:S1:161:ILE:HD12	2.01	0.41
3:S1:157:GLN:O	3:S1:161:ILE:HG13	2.20	0.41
3:S1:178:GLY:O	3:S1:179:SER:OG	5.16	0.41
4:S2:225:LEU:HD22	4:S2:230:TRP:CD1	2.55	0.41
4:S2:53:ILE:HG12	4:S2:72:LEU:HB3	2.03	0.41
5:S3:141:LYS:HB3	5:S3:144:ALA:HA	7.98	0.41
6:S4:58:GLY:HA2	6:S4:61:VAL:HG23	2.26	0.41
7:S5:116:HIS:CD2	27:D5:98:GLN:HB3	3.18	0.41
10:S8:138:ASN:O	10:S8:141:ARG:HB2	2.20	0.41
14:C2:125:ASN:OD1	35:SM:168:UNK:N	4.75	0.41
35:SM:48:ARG:H	35:SM:48:ARG:HG3	1.69	0.41
35:SM:65:THR:HA	35:SM:70:ASN:ND2	2.94	0.41
34:SR:217:ASP:OD1	34:SR:217:ASP:N	2.90	0.41
36:1:1066:G:H2'	36:1:1067:U:C6	2.55	0.41
36:1:1452:A:N3	36:1:2346:C:O2'	2.48	0.41
36:1:1468:A:N6	36:1:1508:C:O2	2.53	0.41
36:1:174:C:H2'	36:1:175:C:C6	2.55	0.41
36:1:2991:A:N3	53:M7:69:ARG:NH2	2.63	0.41
36:1:3137:C:H2'	36:1:3138:U:C6	2.56	0.41
36:1:3222:U:O2'	36:1:3223:A:H5'	2.20	0.41
36:1:981:U:O2'	36:1:982:C:OP1	2.28	0.41
1:2:142:G:C5	1:2:266:A:C6	3.08	0.41
1:2:196:G:O2'	1:2:197:A:P	2.79	0.41
1:2:226:A:H2'	1:2:226:A:N3	2.35	0.41
1:2:281:G:H8	1:2:281:G:OP2	2.02	0.41
1:2:325:G:C6	1:2:344:A:C6	3.09	0.41
36:5:2249:G:H2'	36:5:2250:G:H8	1.84	0.41
65:N9:3:LYS:HD3	36:5:2617:U:H3'	224.38	0.41
52:M6:148:LYS:HE2	36:5:3135:U:OP1	257.57	0.41
36:5:3163:A:C6	36:5:3288:G:C6	3.08	0.41
36:5:315:C:C4	36:5:316:U:C4	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:579:G:H2'	36:5:580:C:C6	2.55	0.41
36:5:742:G:H5'	36:5:743:C:OP1	2.20	0.41
1:6:1111:G:H2'	1:6:1112:G:O4'	2.20	0.41
1:6:1553:G:H2'	1:6:1554:U:H5''	2.01	0.41
1:6:282:C:H2'	1:6:283:U:O4'	2.21	0.41
1:6:837:G:H2'	1:6:838:G:C8	2.55	0.41
37:7:3:U:H2'	37:7:4:U:H6	1.85	0.41
38:8:156:U:H5'	38:8:157:U:OP2	2.20	0.41
13:C1:16:GLN:NE2	13:C1:34:TRP:HB3	3.81	0.41
13:C1:5:LEU:HD13	13:C1:5:LEU:H	5.31	0.41
15:C3:99:ARG:CZ	15:C3:143:SER:HB3	2.50	0.41
17:C5:75:PRO:HG3	17:C5:93:VAL:HG11	2.01	0.41
18:C6:23:LYS:HG2	18:C6:64:ASP:HB2	2.53	0.41
18:C6:7:VAL:HG23	18:C6:22:VAL:O	3.62	0.41
20:C8:115:ARG:O	20:C8:119:ILE:HG12	4.60	0.41
1:2:863:A:O5'	24:D2:57:ARG:HG2	2.21	0.41
25:D3:52:ILE:HG23	25:D3:100:ASP:O	2.20	0.41
1:2:522:U:OP1	26:D4:37:LYS:HB2	2.21	0.41
27:D5:71:ILE:HA	27:D5:71:ILE:HD12	4.90	0.41
28:D6:45:VAL:HB	28:D6:46:GLU:H	1.63	0.41
33:E1:141:CYS:SG	33:E1:143:LYS:HB3	3.49	0.41
40:L3:115:LYS:HA	40:L3:118:PHE:HD1	1.86	0.41
40:L3:230:THR:HB	40:L3:247:ARG:NH1	2.51	0.41
41:L4:152:VAL:CG2	41:L4:172:VAL:HG21	2.50	0.41
42:L5:178:ASN:HA	42:L5:183:TRP:CG	2.64	0.41
43:L6:51:ARG:HD3	43:L6:158:TYR:CZ	2.55	0.41
44:L7:151:ARG:HD2	44:L7:244:ASN:OD1	2.24	0.41
46:L9:117:PHE:HE1	46:L9:178:GLY:HA2	1.91	0.41
47:M0:156:ARG:HG2	47:M0:163:GLN:HG2	2.94	0.41
50:M4:59:ASN:O	50:M4:62:GLN:HG2	4.51	0.41
51:M5:140:LYS:O	51:M5:144:ARG:HG3	2.21	0.41
54:M8:121:CYS:C	54:M8:122:ILE:HD12	2.41	0.41
54:M8:63:SER:O	54:M8:67:ILE:HG13	3.37	0.41
60:N4:38:SER:O	60:N4:42:GLN:HB2	2.84	0.41
36:1:94:G:OP2	64:N8:53:PHE:HD2	2.03	0.41
68:O2:22:SER:HB2	68:O2:30:GLU:HA	2.01	0.41
68:O2:33:ARG:HH11	36:5:944:C:H4'	161.89	0.41
43:L6:6:ALA:HA	68:O2:74:PHE:HE1	1.85	0.41
69:O3:59:VAL:HB	69:O3:60:ARG:H	1.57	0.41
61:N5:45:LYS:HD3	71:O5:75:TYR:HE2	2.73	0.41
71:O5:85:THR:HB	71:O5:88:LEU:HD12	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
71:O5:89:ARG:HD2	38:8:38:U:O4	67.92	0.41
72:O6:51:SER:H	72:O6:54:GLU:HB2	1.85	0.41
74:O8:5:ILE:HD11	74:O8:10:GLN:CD	2.40	0.41
79:Q3:7:LYS:HE3	36:5:1926:C:H2'	253.75	0.41
2:S0:103:THR:HA	2:S0:104:PRO:HD3	1.75	0.41
2:S0:124:THR:O	2:S0:146:LEU:HB2	2.56	0.41
2:S0:31:VAL:N	2:S0:149:LEU:O	2.43	0.41
3:S1:232:HIS:HB3	3:S1:233:GLY:H	2.26	0.41
4:S2:140:ARG:HH22	4:S2:228:ASN:HD21	1.68	0.41
4:S2:63:VAL:HG13	4:S2:68:ILE:HD12	2.03	0.41
6:S4:131:LEU:HD22	6:S4:131:LEU:HA	1.92	0.41
6:S4:141:THR:OG1	6:S4:145:ARG:HB2	2.52	0.41
9:S7:46:ILE:HD13	9:S7:47:ARG:N	2.35	0.41
10:S8:57:ALA:HB1	10:S8:60:ILE:HD12	3.21	0.41
10:S8:78:ILE:HA	10:S8:104:ILE:HG22	2.85	0.41
35:SM:40:PRO:C	35:SM:42:ALA:H	3.00	0.41
36:1:155:G:O2'	72:O6:27:SER:HB3	2.20	0.41
36:1:1675:G:H2'	36:1:1676:A:C8	2.55	0.41
36:1:18:G:OP1	71:O5:81:ARG:NH2	2.53	0.41
36:1:1912:U:C4	36:1:1913:A:C6	3.08	0.41
36:1:2213:A:C6	36:1:2214:A:C6	3.08	0.41
36:1:2263:C:H2'	36:1:2264:U:C6	2.56	0.41
36:1:255:A:C2'	36:1:256:G:H5'	2.50	0.41
36:1:2761:G:N2	36:1:2800:G:O4'	2.54	0.41
36:1:650:C:H2'	36:1:651:G:C8	2.55	0.41
1:2:1046:G:H1	1:2:1072:C:H42	1.69	0.41
1:2:1184:A:O2'	1:2:1209:C:O2'	2.27	0.41
1:2:18:C:H2'	1:2:19:A:C8	2.56	0.41
1:2:610:G:H2'	1:2:614:C:C5	2.56	0.41
1:2:72:A:O2'	1:2:73:U:O4'	2.38	0.41
1:2:782:U:H4'	1:2:783:G:OP2	2.19	0.41
1:2:881:A:H2'	1:2:882:U:O4'	2.21	0.41
37:3:26:C:H2'	37:3:27:A:O4'	2.20	0.41
36:5:1072:G:H2'	36:5:1073:U:C6	2.54	0.41
36:5:1500:G:H2'	36:5:1501:U:O4'	2.20	0.41
36:5:1815:U:O2'	36:5:1816:A:OP2	2.35	0.41
73:O7:9:GLY:HA3	36:5:1852:G:H1'	154.17	0.41
66:O0:50:VAL:HB	36:5:2553:U:O4'	230.02	0.41
79:Q3:62:LYS:NZ	36:5:2554:A:H62	217.24	0.41
78:Q2:39:GLY:N	36:5:2766:U:OP1	170.94	0.41
36:5:982:C:N3	36:5:1102:A:C2	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:1275:A:OP2	1:6:1275:A:H8	2.03	0.41
1:6:1586:A:H2'	1:6:1587:A:O4'	2.20	0.41
1:6:1713:G:C6	1:6:1714:A:C6	3.09	0.41
1:6:329:G:H2'	1:6:330:G:H8	1.86	0.41
1:6:417:A:H5'	1:6:418:G:C5	2.55	0.41
1:6:629:U:H2'	1:6:630:A:H8	1.85	0.41
1:6:722:G:H4'	1:6:723:G:OP1	2.21	0.41
12:C0:44:LYS:HA	12:C0:47:GLN:HB3	2.62	0.41
13:C1:72:THR:HG22	13:C1:124:THR:HG23	2.02	0.41
17:C5:126:VAL:HG21	35:SM:71:ASN:HD21	3.31	0.41
18:C6:11:GLY:O	18:C6:80:ALA:HB1	2.19	0.41
19:C7:81:LYS:HE3	19:C7:81:LYS:HB2	1.81	0.41
20:C8:26:ILE:O	20:C8:31:ALA:HB2	2.31	0.41
24:D2:101:TYR:HB2	24:D2:129:VAL:HG23	3.77	0.41
24:D2:101:TYR:O	24:D2:129:VAL:N	3.01	0.41
25:D3:114:LYS:HB3	25:D3:115:GLY:H	1.67	0.41
20:C8:11:PHE:CD1	27:D5:41:ILE:HG21	4.80	0.41
27:D5:44:GLN:O	27:D5:47:TYR:HB3	2.21	0.41
29:D7:33:LEU:HD13	29:D7:73:LEU:HD21	4.07	0.41
22:D0:65:ILE:HG13	31:D9:43:PHE:CZ	2.55	0.41
32:E0:49:LEU:HD21	32:E0:55:ARG:H	1.86	0.41
39:L2:56:ALA:HB2	39:L2:130:SER:HA	2.71	0.41
40:L3:169:THR:CG2	40:L3:171:LEU:HG	2.50	0.41
40:L3:75:ALA:HB2	36:5:3049:A:C2	245.77	0.41
40:L3:95:THR:OG1	40:L3:98:GLY:O	2.29	0.41
41:L4:187:LEU:HD23	41:L4:187:LEU:HA	1.83	0.41
41:L4:34:ILE:O	41:L4:38:VAL:HG23	2.37	0.41
42:L5:125:VAL:HG11	42:L5:199:ILE:HG21	2.64	0.41
37:3:1:G:C4	42:L5:266:ALA:HA	2.55	0.41
43:L6:69:PHE:N	43:L6:142:ASP:OD2	2.53	0.41
44:L7:103:LEU:HD23	44:L7:103:LEU:HA	1.88	0.41
44:L7:130:ILE:HG21	44:L7:130:ILE:HD13	2.42	0.41
44:L7:184:LEU:HD23	44:L7:184:LEU:HA	2.13	0.41
47:M0:190:VAL:HG22	47:M0:199:PHE:CD1	2.55	0.41
48:M1:14:ILE:HG23	48:M1:129:VAL:HG23	5.71	0.41
48:M1:142:LYS:HD2	48:M1:142:LYS:HA	3.93	0.41
48:M1:92:ARG:HB3	48:M1:173:ASP:OD2	2.20	0.41
51:M5:120:TRP:HE1	51:M5:123:GLN:HB3	2.22	0.41
52:M6:142:SER:C	52:M6:144:SER:H	2.98	0.41
54:M8:54:LEU:HD22	54:M8:58:ASN:HB2	2.03	0.41
57:N1:50:LYS:HB3	57:N1:92:ARG:HH11	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
58:N2:35:LYS:O	58:N2:38:ILE:HG22	2.20	0.41
66:O0:36:GLN:NE2	66:O0:38:LYS:HE2	2.35	0.41
69:O3:38:PRO:HA	69:O3:41:ALA:HB3	3.14	0.41
69:O3:71:VAL:HG13	69:O3:81:VAL:HG11	2.03	0.41
75:O9:28:ARG:HD2	75:O9:36:ARG:HD3	8.56	0.41
77:Q1:15:ARG:HA	77:Q1:18:ARG:HE	1.84	0.41
78:Q2:11:TYR:N	78:Q2:20:HIS:CD2	3.82	0.41
79:Q3:45:LYS:NZ	79:Q3:45:LYS:HB2	2.36	0.41
2:S0:170:ILE:HD12	2:S0:171:GLY:H	1.85	0.41
3:S1:110:LEU:O	3:S1:114:VAL:HG23	2.40	0.41
1:2:6:G:OP2	4:S2:205:ARG:HD2	2.21	0.41
4:S2:115:ILE:HD12	4:S2:212:LYS:HD2	2.03	0.41
5:S3:64:ARG:O	5:S3:66:ILE:N	3.87	0.41
6:S4:191:ARG:HH11	6:S4:245:LYS:HB3	2.23	0.41
6:S4:49:ARG:HH11	6:S4:50:ASN:ND2	2.18	0.41
7:S5:90:ILE:HD13	7:S5:133:VAL:HG11	2.01	0.41
10:S8:60:ILE:HG21	10:S8:179:CYS:HB2	2.02	0.41
11:S9:149:ARG:CZ	1:6:765:G:C6	428.59	0.41
5:S3:125:TYR:CE1	35:SM:134:ASP:OD2	2.74	0.41
34:SR:116:ASP:OD2	34:SR:120:SER:N	2.48	0.41
34:SR:92:TRP:HE3	34:SR:97:GLY:O	2.04	0.41
36:1:115:A:OP1	51:M5:49:ARG:NE	2.54	0.41
36:1:1177:G:N7	69:O3:20:LYS:HD3	2.35	0.41
36:1:1928:G:C6	36:1:1929:G:C4	3.09	0.41
36:1:2353:G:C5	36:1:2354:C:C5	3.09	0.41
36:1:2689:A:H4'	36:1:2690:G:O5'	2.20	0.41
36:1:3113:A:C4	36:1:3114:A:C8	3.09	0.41
36:1:3127:A:H2'	36:1:3128:G:O4'	2.20	0.41
36:1:715:A:H4'	36:1:716:A:OP1	2.20	0.41
1:2:1291:G:H5'	4:S2:119:LYS:NZ	2.36	0.41
1:2:336:G:N2	10:S8:7:SER:HB2	2.35	0.41
1:2:393:C:H41	1:2:400:A:H1'	1.86	0.41
1:2:994:G:C2'	1:2:995:A:H5'	2.50	0.41
38:4:10:A:C5	38:4:11:C:C4	3.08	0.41
62:N6:103:LYS:HE3	36:5:221:A:H61	80.03	0.41
36:5:251:G:C5	36:5:253:A:C6	3.09	0.41
36:5:2726:C:O2	36:5:2728:G:N2	2.53	0.41
36:5:3110:C:H2'	36:5:3111:U:C6	2.56	0.41
36:5:312:C:H1'	36:5:2778:G:N2	2.35	0.41
56:N0:161:LYS:HE3	36:5:3209:A:P	278.17	0.41
36:5:589:A:N6	36:5:610:G:H1'	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:M3:14:PHE:CZ	36:5:665:A:H1'	132.52	0.41
1:6:1163:A:N6	1:6:1164:G:C6	2.89	0.41
1:6:163:G:C6	1:6:164:A:N7	2.88	0.41
1:6:619:A:H5'	1:6:620:A:OP2	2.20	0.41
17:C5:72:LYS:HA	17:C5:73:PRO:HD3	1.77	0.41
21:C9:31:PRO:CG	21:C9:103:LYS:HD3	2.47	0.41
21:C9:38:LYS:HE3	21:C9:38:LYS:HB2	2.27	0.41
21:C9:78:LYS:HE2	1:6:1524:A:H5'	404.09	0.41
23:D1:25:LYS:HB2	23:D1:28:ASP:HB2	5.86	0.41
9:S7:140:VAL:HB	24:D2:52:TYR:HB3	2.14	0.41
24:D2:93:LEU:H	24:D2:93:LEU:HD22	1.86	0.41
25:D3:96:VAL:HA	25:D3:127:VAL:HG21	2.75	0.41
26:D4:20:ARG:C	26:D4:21:LYS:HD2	2.40	0.41
27:D5:41:ILE:HA	27:D5:41:ILE:HD12	1.96	0.41
31:D9:10:HIS:O	31:D9:12:ARG:N	2.53	0.41
31:D9:45:GLU:OE2	1:6:1433:G:N1	407.05	0.41
39:L2:130:SER:HG	39:L2:174:ARG:HH21	2.10	0.41
40:L3:283:TYR:CE1	40:L3:354:VAL:HG11	2.60	0.41
42:L5:11:ALA:O	42:L5:15:ARG:HG3	2.20	0.41
42:L5:285:ARG:O	42:L5:289:LYS:HG3	4.63	0.41
43:L6:136:GLU:O	43:L6:140:VAL:HG23	2.57	0.41
43:L6:30:LEU:HG	43:L6:57:HIS:CD2	2.56	0.41
45:L8:72:PRO:HA	45:L8:73:PRO:HD3	1.92	0.41
48:M1:141:ARG:O	48:M1:145:LYS:NZ	2.49	0.41
48:M1:28:ASP:O	48:M1:32:ARG:HG3	3.54	0.41
48:M1:93:ASP:OD2	48:M1:156:LYS:NZ	3.30	0.41
49:M3:119:TYR:O	49:M3:123:ILE:HG23	2.20	0.41
51:M5:188:ARG:HB2	51:M5:188:ARG:HE	1.65	0.41
53:M7:48:LEU:HB3	53:M7:88:VAL:CG1	2.51	0.41
59:N3:101:VAL:HG21	59:N3:129:VAL:HG22	3.64	0.41
60:N4:6:ASP:HB3	60:N4:11:ALA:H	2.06	0.41
63:N7:104:PRO:O	63:N7:108:GLU:HG3	2.20	0.41
70:O4:16:ARG:HE	70:O4:16:ARG:HB3	3.93	0.41
71:O5:17:LEU:O	71:O5:20:GLN:HB2	2.45	0.41
71:O5:45:LYS:HD2	71:O5:49:LYS:HG3	7.38	0.41
48:M1:60:ARG:NH1	78:Q2:104:LEU:O	3.93	0.41
78:Q2:9:LYS:HE3	78:Q2:22:GLN:CD	4.73	0.41
2:S0:163:ASN:O	2:S0:165:ARG:N	3.20	0.41
3:S1:29:TRP:CD1	3:S1:47:LEU:HG	2.55	0.41
4:S2:179:VAL:HG11	1:6:2:A:H3'	391.03	0.41
5:S3:192:PRO:O	5:S3:195:SER:OG	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S3:195:SER:O	5:S3:197:THR:N	2.54	0.41
6:S4:71:LYS:HA	6:S4:76:VAL:O	2.21	0.41
1:2:186:C:H5''	10:S8:142:LYS:HE2	2.03	0.41
11:S9:129:ILE:HA	11:S9:134:ILE:HG12	3.16	0.41
35:SM:139:GLU:N	35:SM:139:GLU:OE1	2.54	0.41
34:SR:79:TYR:CD1	34:SR:100:TYR:HE1	3.72	0.41
36:1:1018:G:N2	36:1:1019:G:H1'	2.36	0.41
36:1:1541:G:H1'	36:1:1557:A:C5	2.56	0.41
36:1:1636:U:H3	36:1:1710:C:H4'	1.86	0.41
36:1:235:A:H2'	36:1:236:G:O4'	2.21	0.41
36:1:3000:A:H2'	36:1:3001:C:C6	2.56	0.41
36:1:3012:A:N1	36:1:3043:C:C2	2.89	0.41
36:1:35:A:O2'	36:1:36:C:H5'	2.21	0.41
36:1:52:A:N3	36:1:811:U:O2'	2.52	0.41
36:1:607:A:C4	43:L6:26:ARG:NH2	2.88	0.41
36:1:847:A:H2'	36:1:848:A:C8	2.54	0.41
1:2:1795:U:O2	28:D6:10:ARG:HD2	2.21	0.41
1:2:336:G:N2	1:2:338:C:H5'	2.36	0.41
1:2:6:G:H2'	1:2:7:G:C8	2.55	0.41
1:2:993:A:OP1	1:2:1777:G:N2	2.40	0.41
37:3:4:U:H2'	37:3:5:G:C8	2.55	0.41
36:5:1250:G:H2'	36:5:1251:A:C8	2.56	0.41
36:5:1332:A:H2'	36:5:1333:C:C6	2.56	0.41
36:5:2193:U:H5''	36:5:2194:G:H5'	2.01	0.41
36:5:647:A:C2	36:5:2372:A:H2'	2.56	0.41
36:5:2608:G:H2'	36:5:2609:A:H8	1.86	0.41
36:5:2961:G:H2'	36:5:2962:U:C6	2.56	0.41
36:5:3016:A:H2'	36:5:3017:A:C8	2.55	0.41
36:5:3051:U:C2	36:5:3052:G:C8	3.08	0.41
36:5:3291:G:H2'	36:5:3292:A:C8	2.55	0.41
53:M7:118:GLN:HG3	36:5:413:U:O4'	147.46	0.41
43:L6:26:ARG:NH2	36:5:607:A:P	250.06	0.41
36:5:725:G:H5'	36:5:726:G:OP2	2.21	0.41
7:S5:81:ARG:NH2	1:6:1615:C:OP1	364.86	0.41
1:6:620:A:N7	1:6:621:A:C6	2.88	0.41
1:6:828:U:O2'	1:6:829:A:OP1	2.31	0.41
36:5:1055:A:H1'	37:7:81:U:H1'	2.03	0.41
13:C1:71:LEU:HB3	13:C1:88:ARG:NH1	2.43	0.41
14:C2:118:ALA:O	14:C2:120:VAL:N	2.54	0.41
14:C2:81:ASP:OD1	14:C2:85:LYS:HB3	3.73	0.41
19:C7:15:ALA:HA	19:C7:19:ARG:HH21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:C7:41:ILE:O	19:C7:43:SER:N	2.54	0.41
19:C7:95:ARG:N	19:C7:96:SER:HA	4.55	0.41
22:D0:43:LYS:HD2	22:D0:43:LYS:HA	1.89	0.41
28:D6:10:ARG:HD3	28:D6:34:LYS:HA	2.02	0.41
30:D8:15:VAL:HG23	30:D8:27:GLN:O	3.60	0.41
32:E0:31:LYS:HE3	1:6:545:A:OP1	418.76	0.41
40:L3:80:ASP:OD2	40:L3:169:THR:HG23	2.74	0.41
40:L3:199:PHE:C	40:L3:201:LYS:H	2.24	0.41
40:L3:218:ILE:HG12	40:L3:276:THR:HG23	2.03	0.41
41:L4:126:ILE:HD11	41:L4:233:LEU:HD12	3.37	0.41
41:L4:246:ARG:O	41:L4:248:VAL:HG23	3.16	0.41
42:L5:109:THR:OG1	42:L5:110:LEU:N	2.53	0.41
43:L6:170:LYS:HB3	43:L6:172:HIS:CE1	2.68	0.41
44:L7:57:THR:OG1	44:L7:58:ALA:N	3.33	0.41
45:L8:140:VAL:O	45:L8:144:GLU:HG3	2.40	0.41
47:M0:39:LYS:HA	47:M0:86:HIS:CE1	3.20	0.41
47:M0:39:LYS:HG3	47:M0:40:LYS:HG2	4.18	0.41
49:M3:188:ARG:CZ	49:M3:189:GLU:HG3	2.51	0.41
36:1:3178:A:C2	52:M6:115:LYS:HG2	2.55	0.41
54:M8:180:ARG:HA	54:M8:180:ARG:HD2	1.79	0.41
55:M9:167:ARG:HB3	55:M9:167:ARG:NH1	5.12	0.41
55:M9:12:ALA:HB2	55:M9:22:VAL:HG21	3.35	0.41
55:M9:43:LYS:HE3	55:M9:43:LYS:HA	5.28	0.41
55:M9:81:ARG:HG3	55:M9:88:ARG:NH1	2.36	0.41
56:N0:86:GLY:O	56:N0:88:HIS:NE2	4.16	0.41
57:N1:131:GLN:HA	57:N1:132:PRO:HD3	1.86	0.41
57:N1:90:ASN:O	57:N1:91:LEU:HD23	2.21	0.41
61:N5:24:LEU:HD23	61:N5:24:LEU:HA	2.10	0.41
67:O1:53:PRO:O	67:O1:57:GLN:HG3	2.20	0.41
67:O1:84:ASP:OD1	67:O1:84:ASP:N	2.54	0.41
68:O2:8:LYS:HB2	68:O2:8:LYS:HE3	1.82	0.41
49:M3:174:ARG:HB2	72:O6:9:ILE:HD12	2.02	0.41
78:Q2:47:GLN:NE2	78:Q2:53:GLN:OE1	2.53	0.41
3:S1:172:LEU:HA	3:S1:172:LEU:HD23	1.88	0.41
3:S1:33:LYS:HE2	3:S1:41:ARG:HH12	4.14	0.41
4:S2:91:ARG:O	4:S2:91:ARG:NH1	5.68	0.41
5:S3:200:LYS:HB3	5:S3:200:LYS:HE2	1.97	0.41
6:S4:195:ILE:O	6:S4:196:VAL:HB	4.52	0.41
8:S6:58:LYS:HG2	8:S6:105:ASP:O	2.21	0.41
9:S7:39:ARG:NH2	55:M9:185:LEU:HD22	2.36	0.41
10:S8:159:GLN:HG2	10:S8:165:LEU:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:SM:43:ASP:HA	35:SM:44:PRO:HD3	2.54	0.41
34:SR:112:SER:HB2	34:SR:153:GLN:HA	2.18	0.41
34:SR:44:SER:OG	34:SR:59:ARG:HB2	2.21	0.41
36:1:1339:C:H2'	36:1:1340:G:O4'	2.20	0.41
36:1:1713:G:O6	66:O0:28:LYS:NZ	2.30	0.41
36:1:20:A:C6	36:1:21:G:C6	3.09	0.41
36:1:3269:U:H4'	36:1:3270:U:O5'	2.19	0.41
36:1:846:A:H2'	36:1:847:A:O4'	2.20	0.41
1:2:1519:U:H2'	1:2:1520:U:C5	2.55	0.41
1:2:1560:U:H5	1:2:1598:U:HO2'	1.67	0.41
1:2:1660:A:H8	1:2:1660:A:O5'	2.03	0.41
1:2:309:C:H2'	1:2:310:C:C6	2.55	0.41
1:2:540:G:H2'	1:2:540:G:OP2	2.20	0.41
1:2:627:C:H4'	15:C3:117:LEU:CD2	2.51	0.41
1:2:648:G:N1	1:2:649:U:C4	2.89	0.41
36:5:1560:G:O2'	36:5:1561:G:P	2.79	0.41
36:5:2552:C:H5'	36:5:2553:U:OP2	2.20	0.41
40:L3:251:CYS:SG	36:5:2944:U:H1'	224.36	0.41
36:5:297:G:N2	36:5:297:G:OP2	2.47	0.41
36:5:3112:G:C2	36:5:3121:U:C5	3.08	0.41
1:6:1235:C:OP2	1:6:1245:G:H5'	2.21	0.41
1:6:234:G:N3	1:6:234:G:H3'	2.36	0.41
10:S8:31:ARG:O	1:6:331:A:H4'	281.87	0.41
1:6:616:G:C2	1:6:622:A:N7	2.89	0.41
1:6:780:A:OP1	1:6:781:U:H3'	2.21	0.41
1:6:82:U:H2'	1:6:83:G:O4'	2.21	0.41
1:6:836:U:H2'	1:6:837:G:H8	1.86	0.41
1:6:987:G:H5''	1:6:988:A:OP1	2.21	0.41
14:C2:32:LEU:O	14:C2:36:LEU:HB2	2.21	0.41
14:C2:33:ARG:NH2	14:C2:99:GLU:O	8.40	0.41
14:C2:86:VAL:N	14:C2:87:PRO:HD3	2.63	0.41
15:C3:88:LEU:HD23	15:C3:125:LEU:HD13	2.67	0.41
15:C3:27:LYS:H	15:C3:27:LYS:CE	2.29	0.41
15:C3:80:LEU:HA	15:C3:80:LEU:HD22	1.93	0.41
17:C5:31:GLU:O	17:C5:35:LYS:HG2	4.76	0.41
18:C6:113:ASP:OD2	18:C6:114:ARG:N	2.53	0.41
20:C8:145:ARG:HG3	35:SM:68:ARG:CZ	5.62	0.41
21:C9:134:ARG:O	21:C9:138:GLN:HG3	2.21	0.41
21:C9:28:LEU:O	21:C9:29:GLU:HB2	2.19	0.41
22:D0:106:ILE:O	22:D0:107:THR:OG1	2.30	0.41
23:D1:71:ARG:NE	29:D7:4:VAL:HG11	3.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:D3:65:ASN:HD22	25:D3:65:ASN:HA	1.56	0.41
40:L3:100:ARG:NH2	36:5:3244:A:OP 2	250.27	0.41
40:L3:287:LYS:HD2	40:L3:287:LYS:HA	4.80	0.41
41:L4:333:VAL:HG23	41:L4:337:GLU:HG3	3.31	0.41
41:L4:81:GLY:HA3	36:5:357:A:O4'	129.23	0.41
42:L5:279:LYS:HD3	42:L5:279:LYS:O	2.21	0.41
42:L5:46:THR:HA	42:L5:47:PRO:HD3	2.34	0.41
36:1:591:G:H1'	43:L6:19:LYS:HG3	2.03	0.41
47:M0:161:GLY:O	47:M0:163:GLN:NE2	2.60	0.41
47:M0:184:LYS:HE3	47:M0:189:GLU:OE1	2.21	0.41
47:M0:210:ILE:HG23	47:M0:217:PHE:CE2	2.56	0.41
49:M3:2:ALA:N	64:N8:31:GLY:O	2.57	0.41
50:M4:113:THR:H	50:M4:116:GLU:HB3	3.96	0.41
50:M4:34:ALA:HB2	50:M4:85:TRP:HZ3	2.04	0.41
50:M4:70:PHE:CE2	50:M4:72:LEU:HD23	2.56	0.41
52:M6:62:THR:HG22	52:M6:65:ASN:H	1.94	0.41
52:M6:35:VAL:HG21	52:M6:80:PHE:HE2	1.86	0.41
53:M7:119:VAL:HA	53:M7:145:HIS:O	2.26	0.41
53:M7:128:ARG:HG2	53:M7:136:ILE:HG21	5.51	0.41
57:N1:102:ARG:HD3	57:N1:102:ARG:HA	1.75	0.41
36:1:2698:G:O2'	57:N1:12:ARG:HG2	2.20	0.41
44:L7:77:VAL:CG2	57:N1:139:ARG:HG2	2.51	0.41
58:N2:43:VAL:HG21	58:N2:50:LEU:HA	2.02	0.41
63:N7:3:LYS:O	63:N7:6:LYS:HG3	2.21	0.41
63:N7:53:VAL:HA	63:N7:57:HIS:CD2	2.56	0.41
64:N8:16:SER:O	64:N8:17:ALA:HB3	2.21	0.41
65:N9:38:LYS:HG3	36:5:2738:A:H5"	213.03	0.41
65:N9:53:ALA:O	65:N9:56:ALA:HB3	2.74	0.41
68:O2:67:SER:HB3	68:O2:68:PRO:HD2	2.02	0.41
36:1:3173:G:C2	69:O3:96:ALA:HB2	2.56	0.41
72:O6:61:ILE:HA	72:O6:65:GLY:O	2.20	0.41
72:O6:69:ALA:O	72:O6:73:ALA:N	2.53	0.41
72:O6:97:SER:HB3	72:O6:98:ARG:H	1.65	0.41
73:O7:19:CYS:HB2	73:O7:27:PHE:HB2	2.02	0.41
73:O7:75:LYS:HE3	73:O7:75:LYS:HB3	1.91	0.41
78:Q2:14:GLY:O	78:Q2:18:ARG:HG3	2.20	0.41
2:S0:160:ILE:HA	2:S0:161:PRO:HD2	1.90	0.41
2:S0:62:ARG:HD3	23:D1:36:VAL:HG13	2.03	0.41
3:S1:128:LYS:HG2	3:S1:129:THR:O	2.21	0.41
3:S1:180:THR:HG23	3:S1:183:GLN:HE21	9.87	0.41
3:S1:229:MET:O	3:S1:231:LEU:N	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S1:35:PRO:HG3	3:S1:98:THR:O	2.20	0.41
4:S2:129:ILE:HG22	4:S2:133:LYS:HE3	2.67	0.41
6:S4:12:LEU:HD13	6:S4:12:LEU:HA	1.89	0.41
6:S4:31:PRO:HD2	6:S4:38:LEU:HD13	2.92	0.41
6:S4:46:VAL:O	6:S4:50:ASN:HB2	2.32	0.41
9:S7:25:VAL:O	9:S7:28:GLU:HB2	3.38	0.41
1:2:856:A:N6	9:S7:96:ARG:HB3	2.35	0.41
9:S7:9:LEU:HD11	9:S7:17:GLU:HB3	4.22	0.41
10:S8:110:ARG:HH12	10:S8:160:PHE:HB3	3.10	0.41
11:S9:99:LEU:HD12	11:S9:100:LYS:H	1.85	0.41
35:SM:121:LYS:HD3	35:SM:121:LYS:HA	1.95	0.41
34:SR:93:ASP:HB2	34:SR:100:TYR:CE1	2.56	0.41
34:SR:121:MET:SD	34:SR:183:LEU:HD23	6.26	0.41
34:SR:172:ALA:HB2	34:SR:202:LEU:HD22	2.62	0.41
36:1:945:C:H1'	36:1:1407:A:H1'	2.03	0.41
36:1:1719:G:H5''	55:M9:110:ARG:HH22	1.86	0.41
36:1:1801:U:O5'	36:1:1801:U:H6	2.04	0.41
36:1:2367:A:H2'	36:1:2368:A:C8	2.56	0.41
36:1:2651:G:H5''	36:1:2652:U:O4'	2.21	0.41
36:1:2855:U:H2'	36:1:2856:G:O4'	2.21	0.41
36:1:3164:C:O2'	36:1:3165:A:H8	2.04	0.41
36:1:3279:A:N7	69:O3:54:ARG:NH2	2.68	0.41
36:1:3297:U:O4	40:L3:124:LYS:NZ	2.54	0.41
36:1:1940:G:N2	36:1:3362:A:C8	2.86	0.41
36:1:3373:U:OP2	67:O1:102:LYS:HE3	2.20	0.41
36:1:792:G:H2'	36:1:793:C:C6	2.56	0.41
36:1:924:G:N3	36:1:2414:G:O2'	2.52	0.41
1:2:936:G:OP1	1:2:1075:C:H1'	2.21	0.41
1:2:1391:A:H61	1:2:1407:U:H3	1.69	0.41
1:2:1471:A:N3	1:2:1474:G:O2'	2.40	0.41
1:2:1753:A:N1	1:2:1754:A:N6	2.68	0.41
1:2:1774:G:H2'	1:2:1775:U:O4'	2.21	0.41
1:2:823:G:O2'	1:2:824:G:O4'	2.31	0.41
1:2:940:A:H2'	1:2:941:A:O4'	2.21	0.41
37:3:90:U:C4	37:3:91:G:C5	3.09	0.41
38:4:154:C:H2'	38:4:155:A:O4'	2.20	0.41
36:5:1078:U:H1'	36:5:1082:U:O2	2.21	0.41
36:5:1288:U:H2'	36:5:1289:G:C8	2.56	0.41
36:5:210:U:O2'	36:5:229:G:O2'	2.37	0.41
36:5:2833:A:C2	36:5:2834:G:C8	3.09	0.41
36:5:3275:U:H4'	36:5:3276:G:OP2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:5:529:A:H2'	36:5:530:G:O4'	2.20	0.41
1:6:1120:U:H3	1:6:1127:G:H1	1.69	0.41
1:6:1255:G:H4'	1:6:1256:A:OP1	2.21	0.41
1:6:1614:A:O2'	1:6:1615:C:H5'	2.20	0.41
25:D3:140:LYS:NZ	1:6:31:C:OP1	378.46	0.41
1:6:44:U:H2'	1:6:45:U:C5	2.56	0.41
1:6:706:A:H2'	1:6:707:A:O4'	2.20	0.41
55:M9:163:ARG:HD3	1:6:813:U:C2	304.21	0.41
38:8:81:U:H1'	38:8:82:U:H6	1.85	0.41
12:C0:55:VAL:HB	12:C0:67:THR:O	2.21	0.41
14:C2:54:ARG:O	14:C2:56:GLU:N	2.54	0.41
15:C3:20:ARG:HD3	24:D2:56:HIS:CD2	5.04	0.41
15:C3:4:MET:HG3	15:C3:5:HIS:H	1.85	0.41
19:C7:29:GLN:HB3	34:SR:85:TRP:HZ3	1.86	0.41
5:S3:209:ILE:HB	19:C7:38:ILE:O	3.68	0.41
20:C8:113:LEU:O	20:C8:117:LYS:HG3	2.60	0.41
21:C9:14:PHE:CE1	21:C9:135:ILE:HD11	2.56	0.41
21:C9:30:VAL:HA	21:C9:31:PRO:HD2	1.80	0.41
21:C9:14:PHE:CE2	21:C9:63:ARG:HB2	2.56	0.41
25:D3:10:ASN:O	1:6:632:U:H5''	333.02	0.41
25:D3:72:VAL:O	25:D3:84:THR:HA	2.62	0.41
7:S5:119:ASP:HB3	27:D5:100:ILE:HD13	2.17	0.41
7:S5:189:THR:HG21	27:D5:98:GLN:HE22	4.12	0.41
28:D6:31:PRO:O	28:D6:34:LYS:N	3.00	0.41
39:L2:129:ALA:HB3	39:L2:132:ASN:OD1	2.21	0.41
36:1:2245:C:H4'	39:L2:221:LYS:O	2.21	0.41
40:L3:296:THR:HG21	40:L3:356:LEU:C	2.61	0.41
41:L4:118:LYS:O	41:L4:121:ALA:HB3	2.84	0.41
41:L4:166:VAL:O	41:L4:170:LYS:HG3	2.21	0.41
41:L4:181:VAL:HG13	41:L4:201:GLN:HB2	2.02	0.41
41:L4:198:ARG:HD3	41:L4:199:TRP:NE1	2.36	0.41
41:L4:258:LEU:HD12	41:L4:258:LEU:HA	1.87	0.41
42:L5:218:ARG:HA	42:L5:221:GLU:CD	2.41	0.41
42:L5:95:TRP:CD1	42:L5:198:TYR:HB3	4.43	0.41
44:L7:156:ILE:HD12	44:L7:161:VAL:HB	2.02	0.41
45:L8:49:TYR:O	36:5:2523:A:H2'	170.76	0.41
46:L9:106:LYS:H	46:L9:109:ALA:CB	2.34	0.41
48:M1:100:GLY:HA3	48:M1:154:THR:HB	2.37	0.41
48:M1:131:MET:HB3	48:M1:131:MET:HE3	2.37	0.41
49:M3:60:ALA:HB3	49:M3:65:TYR:O	2.20	0.41
36:1:3229:G:O5'	50:M4:137:LYS:NZ	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:M7:21:TYR:H	53:M7:145:HIS:CE1	2.38	0.41
58:N2:30:PRO:HA	58:N2:33:TYR:HB3	2.03	0.41
59:N3:74:MET:HE3	59:N3:74:MET:HB3	4.78	0.41
63:N7:17:ARG:HB2	36:5:1635:G:O6	202.17	0.41
63:N7:24:VAL:HG11	63:N7:87:LEU:HB3	2.03	0.41
54:M8:157:PRO:HD3	64:N8:47:LYS:HB2	2.02	0.41
66:O0:30:THR:HB	66:O0:91:SER:HB2	2.13	0.41
63:N7:136:PHE:CD2	70:O4:89:ILE:HG13	2.55	0.41
74:O8:27:ILE:HG22	74:O8:41:THR:HG22	2.03	0.41
75:O9:27:ILE:C	75:O9:29:LEU:H	2.79	0.41
78:Q2:4:VAL:O	78:Q2:94:GLY:N	4.55	0.41
79:Q3:83:ILE:HD13	79:Q3:83:ILE:HA	1.86	0.41
2:S0:52:LYS:HG2	2:S0:52:LYS:H	1.47	0.41
3:S1:47:LEU:HD12	3:S1:47:LEU:H	2.57	0.41
5:S3:134:CYS:N	5:S3:157:LEU:HD11	2.36	0.41
5:S3:194:LYS:O	5:S3:196:ARG:N	2.44	0.41
6:S4:52:LEU:HB3	6:S4:54:TYR:CD2	2.61	0.41
7:S5:203:LYS:HE2	7:S5:203:LYS:HB2	4.31	0.41
8:S6:173:PRO:HB2	8:S6:174:LYS:H	1.53	0.41
10:S8:67:TRP:CZ2	10:S8:156:VAL:HG11	2.55	0.41
11:S9:127:VAL:HG12	11:S9:131:GLN:OE1	2.21	0.41
1:2:380:U:H5	11:S9:5:PRO:HB3	1.86	0.41
19:C7:29:GLN:HB2	34:SR:38:ARG:HH12	4.08	0.41
34:SR:67:ILE:HD12	34:SR:85:TRP:CZ3	2.64	0.41
36:1:1217:A:C2	36:1:1289:G:C4	3.09	0.41
36:1:1286:A:N3	36:1:1287:A:H1'	2.36	0.41
36:1:1350:A:C2'	36:1:1351:U:H5'	2.51	0.41
36:1:1668:G:C5	36:1:1669:C:C5	3.09	0.41
36:1:1830:G:O3'	61:N5:91:ASN:HB2	2.21	0.41
36:1:2376:G:C6	36:1:2377:G:C6	3.08	0.41
36:1:2620:G:H1	84:C:74:CH:HN3	1.68	0.41
36:1:2374:C:C4	36:1:2941:A:C4	3.09	0.41
36:1:3059:G:OP2	67:O1:69:TYR:OH	2.37	0.41
36:1:353:G:O6	73:O7:52:LYS:NZ	2.52	0.41
36:1:506:U:H2'	36:1:507:U:O4'	2.21	0.41
36:1:535:G:N1	36:1:555:U:C2	2.89	0.41
36:1:547:G:H4'	36:1:548:G:OP1	2.21	0.41
36:1:602:A:H2'	36:1:603:A:C8	2.56	0.41
36:1:733:G:N2	36:1:736:A:OP2	2.50	0.41
36:1:980:A:N7	36:1:981:U:C4	2.89	0.41
1:2:1116:A:H1'	1:2:1131:A:N6	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1256:A:H4'	1:2:1257:U:O5'	2.20	0.41
1:2:1301:U:OP1	4:S2:88:LYS:HB2	2.21	0.41
1:2:1486:G:H1'	1:2:1592:A:O2'	2.21	0.41
1:2:315:A:N3	1:2:316:A:H1'	2.36	0.41
1:2:461:G:OP1	11:S9:2:PRO:HG2	2.21	0.41
37:3:23:A:C6	37:3:24:A:C6	3.09	0.41
37:3:30:G:C6	37:3:31:U:C4	3.09	0.41
67:O1:57:GLN:HG2	36:5:1475:A:H4'	147.03	0.41
36:5:1701:C:H2'	36:5:1702:U:O4'	2.21	0.41
36:5:1767:C:H2'	36:5:1768:U:C6	2.56	0.41
36:5:1816:A:O2'	36:5:1817:G:OP1	2.36	0.41
55:M9:104:ARG:HH12	36:5:1949:G:H5''	215.22	0.41
36:5:260:C:H2'	36:5:261:U:H6	1.85	0.41
52:M6:117:ARG:NH2	36:5:3180:A:O2'	276.87	0.41
36:5:3274:A:O2'	36:5:3275:U:OP1	2.37	0.41
36:5:547:G:C6	36:5:548:G:H1'	2.56	0.41
51:M5:81:TYR:OH	36:5:908:G:H3'	165.31	0.41
36:5:952:A:O3'	36:5:968:G:N2	2.54	0.41
1:6:139:C:H4'	1:6:140:A:H5'	2.03	0.41
1:6:315:A:H4'	1:6:316:A:H4'	2.01	0.41
15:C3:45:LEU:HD23	15:C3:45:LEU:HA	4.15	0.41
16:C4:13:VAL:HG13	16:C4:77:THR:H	1.86	0.41
19:C7:71:PHE:C	19:C7:73:LEU:H	2.23	0.41
1:2:1566:U:H5''	20:C8:39:GLY:N	2.36	0.41
23:D1:53:TYR:OH	23:D1:76:ASP:HB2	3.11	0.41
24:D2:49:GLU:O	24:D2:64:GLN:HG2	2.74	0.41
24:D2:26:LEU:HD23	24:D2:62:VAL:HG22	2.80	0.41
24:D2:23:ARG:HH11	24:D2:66:ASN:HA	1.85	0.41
25:D3:107:PHE:CE1	25:D3:114:LYS:HD3	5.57	0.41
39:L2:181:LYS:HB2	36:5:860:G:C5	211.86	0.41
39:L2:47:GLN:CD	39:L2:60:LYS:HD2	5.90	0.41
39:L2:70:ARG:NH2	36:5:2522:G:O6	175.84	0.41
39:L2:79:ASN:H	39:L2:82:VAL:HG11	2.79	0.41
40:L3:313:HIS:O	40:L3:333:LYS:HD2	2.21	0.41
46:L9:27:VAL:HG12	46:L9:82:VAL:HG11	2.03	0.41
48:M1:79:ILE:HG22	48:M1:127:PHE:HE2	1.86	0.41
35:SM:39:PRO:HD3	48:M1:52:TYR:CE1	3.10	0.41
49:M3:31:LYS:O	49:M3:35:ARG:HG3	4.42	0.41
49:M3:67:ARG:HG3	49:M3:67:ARG:H	1.52	0.41
51:M5:21:PHE:CD2	51:M5:22:LEU:HD13	2.56	0.41
52:M6:73:PHE:HB3	52:M6:78:ARG:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:M6:85:ARG:HD3	52:M6:90:HIS:ND1	2.41	0.41
53:M7:39:TRP:O	53:M7:114:VAL:HG12	2.24	0.41
57:N1:62:GLY:HA3	57:N1:76:ILE:HD13	2.03	0.41
58:N2:85:LYS:HG2	36:5:1682:U:C6	152.14	0.41
36:1:3051:U:OP1	60:N4:17:ARG:HD3	2.21	0.41
63:N7:24:VAL:HG23	63:N7:44:ALA:O	2.25	0.41
63:N7:77:TYR:HB3	66:O0:35:ARG:CZ	3.10	0.41
64:N8:81:LEU:HD21	64:N8:109:TYR:HE1	1.85	0.41
66:O0:16:LEU:HD12	66:O0:97:ASP:HB3	2.03	0.41
70:O4:99:LYS:HB2	70:O4:99:LYS:HE3	1.73	0.41
71:O5:47:VAL:HA	71:O5:50:SER:HB3	2.03	0.41
71:O5:93:THR:H	71:O5:96:GLU:HG3	1.86	0.41
73:O7:52:LYS:O	73:O7:55:ARG:HB3	2.20	0.41
78:Q2:25:VAL:HG12	78:Q2:93:LEU:HD12	2.02	0.41
79:Q3:22:LEU:HA	79:Q3:22:LEU:HD23	1.93	0.41
4:S2:168:ARG:HD2	1:6:1097:U:H6	380.73	0.41
5:S3:115:ILE:HD11	5:S3:138:VAL:HG21	2.03	0.41
5:S3:207:THR:HG1	19:C7:40:THR:HG1	3.73	0.41
5:S3:53:THR:O	5:S3:90:ARG:NH2	2.53	0.41
6:S4:104:ASP:N	6:S4:108:ARG:O	4.33	0.41
7:S5:113:ILE:HD13	7:S5:190:ILE:HB	2.03	0.41
9:S7:9:LEU:HD23	9:S7:10:SER:H	4.59	0.41
11:S9:121:SER:HB3	11:S9:124:HIS:HB2	4.61	0.41
34:SR:42:LEU:HB2	34:SR:61:PHE:CD2	3.38	0.41
36:1:1422:G:H21	43:L6:5:LYS:NZ	2.19	0.41
36:1:1562:C:H4'	36:1:1563:C:OP1	2.21	0.41
36:1:1595:U:C2	36:1:1596:C:C5	3.09	0.41
36:1:2822:U:H2'	36:1:2823:G:O4'	2.21	0.41
36:1:2824:G:H2'	36:1:2825:C:C6	2.55	0.41
36:1:3215:A:H62	50:M4:122:VAL:HG13	1.85	0.41
36:1:855:U:H2'	36:1:856:G:O4'	2.21	0.41
36:1:966:U:C2	36:1:967:A:C8	3.09	0.41
1:2:1081:A:H2	1:2:1082:C:H41	1.68	0.41
1:2:1603:U:H2'	1:2:1604:U:C6	2.56	0.41
1:2:1155:G:N2	1:2:1624:C:C2	2.88	0.41
1:2:1149:G:H1'	1:2:1765:A:C4	2.56	0.41
1:2:316:A:H2'	1:2:317:C:H6	1.86	0.41
1:2:452:A:H3'	1:2:453:U:C6	2.56	0.41
1:2:55:A:OP1	26:D4:112:LYS:NZ	2.41	0.41
1:2:627:C:H2'	1:2:628:G:O4'	2.21	0.41
1:2:720:G:H2'	1:2:720:G:N3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:3:98:C:N3	37:3:99:G:C8	2.89	0.41
36:5:1366:A:H2'	36:5:1366:A:N3	2.36	0.41
36:5:1560:G:H1	36:5:1579:C:H42	1.69	0.41
36:5:1622:U:H2'	36:5:1623:G:C8	2.56	0.41
36:5:1648:A:H2'	36:5:1649:U:O4'	2.20	0.41
36:5:1691:U:H2'	36:5:1692:U:C6	2.55	0.41
36:5:1916:U:H2'	36:5:1917:C:H6	1.86	0.41
36:5:2097:U:H2'	36:5:2098:C:C6	2.56	0.41
36:5:2258:U:H5''	36:5:2259:A:OP2	2.21	0.41
36:5:3014:U:H2'	36:5:3015:G:H8	1.85	0.41
36:5:3205:G:H2'	36:5:3206:C:C5	2.56	0.41
36:5:3287:U:N3	36:5:3288:G:N7	2.69	0.41
36:5:3163:A:C6	36:5:3288:G:O6	2.74	0.41
36:5:3302:U:H3	36:5:3312:U:H3	1.67	0.41
36:5:707:U:H2'	36:5:708:G:C8	2.56	0.41
36:5:830:A:H4'	36:5:1866:C:C4	2.55	0.41
1:6:1662:G:H2'	1:6:1663:G:H8	1.86	0.41
10:S8:24:LYS:O	1:6:400:A:H5''	307.42	0.41
1:6:447:U:C4	1:6:448:C:C4	3.09	0.41
8:S6:175:ILE:HG12	1:6:78:A:H1'	337.89	0.41
1:6:825:U:H6	1:6:825:U:OP2	2.04	0.41
15:C3:128:TYR:CE1	1:6:964:U:H5''	322.36	0.41
13:C1:112:SER:C	13:C1:114:ALA:H	2.24	0.41
15:C3:4:MET:O	15:C3:5:HIS:ND1	6.03	0.41
18:C6:37:THR:HA	18:C6:49:TYR:OH	2.64	0.41
19:C7:27:ASP:HB3	19:C7:30:THR:HG23	2.03	0.41
20:C8:28:ILE:HG22	20:C8:56:LYS:O	2.21	0.41
2:S0:55:GLU:HG2	23:D1:79:LEU:HD22	6.86	0.41
13:C1:99:ARG:HD3	25:D3:8:GLY:O	2.24	0.41
26:D4:18:LEU:HA	26:D4:18:LEU:HD23	1.88	0.41
28:D6:62:TYR:CG	28:D6:63:ALA:N	2.87	0.41
28:D6:71:LEU:HD13	28:D6:73:TYR:OH	3.93	0.41
30:D8:30:VAL:HB	30:D8:31:GLU:H	2.51	0.41
39:L2:20:THR:O	39:L2:23:ARG:HG3	2.20	0.41
36:1:3045:G:O3'	40:L3:275:ARG:NH1	2.54	0.41
40:L3:46:PHE:CZ	40:L3:84:VAL:HB	2.56	0.41
41:L4:131:VAL:O	41:L4:135:VAL:HG23	2.75	0.41
42:L5:257:GLU:O	42:L5:258:LYS:HD3	5.16	0.41
42:L5:278:SER:H	42:L5:281:GLU:CD	3.22	0.41
42:L5:69:ILE:HD13	57:N1:28:SER:HB2	2.03	0.41
41:L4:349:THR:HG21	44:L7:64:GLN:HE21	5.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:L7:88:ARG:CZ	44:L7:103:LEU:HD13	2.50	0.41
45:L8:140:VAL:HG21	51:M5:3:ALA:HB2	2.03	0.41
46:L9:22:SER:OG	46:L9:23:ARG:N	2.53	0.41
47:M0:191:LYS:HE3	47:M0:191:LYS:HB3	4.34	0.41
49:M3:176:GLU:HA	49:M3:183:ARG:HH12	1.86	0.41
49:M3:80:VAL:HG22	49:M3:85:LEU:O	2.20	0.41
51:M5:118:SER:HB3	51:M5:132:VAL:HG13	2.29	0.41
53:M7:25:SER:CB	53:M7:28:ASN:HB2	2.79	0.41
53:M7:39:TRP:HB2	53:M7:44:ALA:HB2	2.27	0.41
53:M7:46:LYS:O	53:M7:50:GLN:HG3	2.20	0.41
55:M9:35:ALA:O	55:M9:37:SER:N	3.00	0.41
57:N1:12:ARG:HD2	57:N1:13:TYR:HE1	1.81	0.41
62:N6:100:HIS:ND1	62:N6:102:SER:OG	2.46	0.41
62:N6:5:SER:C	62:N6:7:ASP:N	3.05	0.41
63:N7:105:SER:HA	63:N7:108:GLU:OE1	2.22	0.41
63:N7:22:LYS:HE3	63:N7:134:LEU:HB2	2.84	0.41
63:N7:4:PHE:HE1	63:N7:82:PRO:HG3	1.85	0.41
49:M3:153:ASP:OD2	64:N8:101:VAL:HG11	4.70	0.41
65:N9:43:HIS:O	65:N9:47:LEU:HG	2.21	0.41
66:O0:104:LEU:HD12	66:O0:105:ALA:N	2.65	0.41
68:O2:71:HIS:ND1	68:O2:118:LYS:HD3	2.35	0.41
68:O2:7:PRO:HG2	68:O2:63:THR:HG23	3.79	0.41
69:O3:67:MET:HE1	69:O3:87:ASN:O	5.61	0.41
78:Q2:80:ARG:HB2	36:5:2769:A:O3'	195.38	0.41
2:S0:142:PRO:HB3	23:D1:34:ILE:HD13	2.12	0.41
3:S1:62:LYS:O	3:S1:88:VAL:HB	2.21	0.41
4:S2:101:VAL:HG22	4:S2:115:ILE:HG23	2.03	0.41
5:S3:64:ARG:NH1	5:S3:68:GLU:OE1	3.24	0.41
6:S4:112:HIS:NE2	6:S4:237:SER:O	2.54	0.41
6:S4:29:PRO:O	6:S4:30:ARG:HB3	4.58	0.41
6:S4:87:MET:SD	6:S4:123:LEU:HB2	2.77	0.41
7:S5:22:PRO:HA	7:S5:34:GLN:OE1	6.23	0.41
1:2:1473:U:C5	7:S5:98:MET:HA	2.51	0.41
8:S6:39:GLU:HA	8:S6:42:GLY:O	2.21	0.41
9:S7:75:THR:HG22	9:S7:161:GLN:NE2	4.72	0.41
10:S8:74:LYS:HG2	10:S8:108:PRO:HB2	2.03	0.41
11:S9:112:GLN:O	11:S9:116:LEU:HB3	4.75	0.41
5:S3:115:ILE:HG21	35:SM:110:TRP:HA	2.02	0.41
35:SM:58:GLU:O	35:SM:62:ARG:HB2	2.78	0.41
34:SR:244:ALA:O	34:SR:252:LEU:HD12	3.35	0.41
34:SR:305:TYR:CE2	34:SR:311:ARG:HD2	3.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:139:G:H2'	36:1:140:C:C6	2.55	0.40
36:1:1580:A:H1'	36:1:1581:C:H5	1.86	0.40
36:1:1793:C:O2	39:L2:188:LYS:HE3	2.21	0.40
36:1:1898:G:H2'	36:1:1899:G:O4'	2.20	0.40
36:1:2101:C:H1'	36:1:2102:U:O5'	2.22	0.40
36:1:2240:G:H2'	36:1:2241:U:O4'	2.21	0.40
36:1:256:G:H4'	71:O5:111:PHE:HZ	1.86	0.40
36:1:776:U:C5	36:1:2719:U:O2	2.74	0.40
36:1:2747:A:H4'	42:L5:174:PRO:O	2.21	0.40
36:1:2878:G:H2'	36:1:2879:C:H6	1.85	0.40
36:1:36:C:H2'	36:1:37:U:H5'	2.03	0.40
1:2:1107:G:C6	1:2:1108:G:C6	3.09	0.40
1:2:1796:C:H1'	28:D6:7:SER:OG	2.22	0.40
1:2:182:A:H2'	1:2:183:U:C6	2.56	0.40
1:2:377:G:H4'	1:2:379:U:O4	2.22	0.40
1:2:514:G:O2'	1:2:515:A:H5'	2.20	0.40
1:2:707:A:H61	1:2:730:G:H22	1.69	0.40
36:5:1008:U:H2'	36:5:1009:A:O4'	2.21	0.40
36:5:1077:U:H2'	36:5:1078:U:C6	2.57	0.40
36:5:1341:U:H2'	36:5:1342:C:C6	2.56	0.40
36:5:1911:A:H2	36:5:2122:G:C8	2.40	0.40
36:5:2423:U:H2'	36:5:2424:A:C8	2.56	0.40
78:Q2:9:LYS:O	36:5:2713:U:H3'	223.78	0.40
36:5:3241:G:H2'	36:5:3245:A:N7	2.37	0.40
36:5:3362:A:C2	36:5:3363:U:C2	3.09	0.40
36:5:359:U:H4'	36:5:817:A:N6	2.36	0.40
36:5:661:G:C5	36:5:802:C:C6	3.09	0.40
41:L4:119:ARG:NH2	36:5:696:C:OP2	103.02	0.40
1:6:1177:C:H4'	1:6:1189:A:N1	2.36	0.40
1:6:1241:G:C4	1:6:1242:A:C8	3.10	0.40
1:6:445:A:N1	1:6:462:G:O2'	2.46	0.40
1:6:473:A:H4'	1:6:768:C:O2	2.20	0.40
1:6:934:C:C4	1:6:1077:C:H4'	2.55	0.40
37:7:113:C:C4	37:7:114:U:C4	3.09	0.40
38:8:25:G:C6	38:8:26:U:C2	3.08	0.40
38:8:68:G:C5	38:8:69:U:C5	3.09	0.40
12:C0:44:LYS:HD3	12:C0:44:LYS:HA	1.72	0.40
13:C1:45:PRO:O	13:C1:49:ILE:HG13	3.43	0.40
14:C2:43:ARG:HD3	14:C2:43:ARG:H	1.86	0.40
14:C2:66:VAL:HG23	14:C2:69:ALA:HA	3.64	0.40
1:2:1180:C:O2'	17:C5:128:HIS:ND1	2.48	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:C5:130:ARG:HD3	17:C5:130:ARG:HA	1.71	0.40
17:C5:53:PRO:O	17:C5:57:MET:HB2	3.80	0.40
20:C8:134:ARG:HG3	1:6:1545:A:OP2	356.06	0.40
23:D1:59:VAL:HG23	23:D1:59:VAL:H	1.65	0.40
12:C0:61:TRP:HZ3	31:D9:22:ARG:HD3	1.86	0.40
40:L3:41:VAL:HG22	40:L3:186:GLY:H	1.86	0.40
40:L3:285:VAL:HG13	40:L3:322:ILE:CD1	2.52	0.40
40:L3:285:VAL:HG13	40:L3:322:ILE:HD13	2.02	0.40
40:L3:29:VAL:HG22	40:L3:218:ILE:HD13	3.56	0.40
41:L4:303:GLY:H	36:5:1347:U:H5"	197.26	0.40
42:L5:219:PHE:HE1	42:L5:227:LEU:HD11	3.33	0.40
42:L5:261:THR:H	42:L5:264:GLN:HG3	1.86	0.40
36:1:3271:G:C5	43:L6:108:LYS:HE3	2.56	0.40
45:L8:146:LYS:HE3	45:L8:173:MET:HE3	2.02	0.40
45:L8:147:LYS:HE2	45:L8:147:LYS:HB3	1.82	0.40
47:M0:74:LYS:HB2	47:M0:74:LYS:HE3	1.83	0.40
48:M1:166:LYS:HE2	48:M1:166:LYS:HB2	2.12	0.40
48:M1:16:LYS:HE3	48:M1:130:VAL:HG11	2.03	0.40
48:M1:52:TYR:HA	48:M1:61:ARG:HG3	2.02	0.40
51:M5:59:PHE:CE2	51:M5:142:ILE:HD11	4.23	0.40
52:M6:65:ASN:HB3	52:M6:68:ARG:CD	2.51	0.40
52:M6:73:PHE:CE1	36:5:3007:U:H5'	242.87	0.40
54:M8:148:GLU:OE1	54:M8:151:ARG:NE	3.60	0.40
36:1:2700:G:OP1	57:N1:17:ARG:HB2	2.21	0.40
62:N6:27:ARG:HA	62:N6:30:LEU:HB2	2.03	0.40
62:N6:30:LEU:HD23	62:N6:30:LEU:HA	1.91	0.40
68:O2:38:ILE:H	68:O2:38:ILE:HG12	1.98	0.40
69:O3:6:ARG:HG3	69:O3:8:TYR:H	2.01	0.40
72:O6:36:ARG:O	72:O6:40:VAL:HG23	2.20	0.40
2:S0:76:ILE:O	2:S0:124:THR:HG23	2.21	0.40
2:S0:53:THR:O	2:S0:57:LEU:HB2	2.57	0.40
3:S1:71:ALA:HB2	3:S1:79:HIS:O	2.20	0.40
4:S2:115:ILE:HD13	4:S2:208:GLU:HG2	2.03	0.40
5:S3:176:LEU:HA	5:S3:181:VAL:HG12	2.03	0.40
5:S3:76:ARG:HG3	12:C0:65:TYR:HE1	1.86	0.40
6:S4:121:TYR:CD2	6:S4:161:LYS:HE3	2.57	0.40
6:S4:213:SER:O	6:S4:213:SER:OG	2.78	0.40
6:S4:251:GLU:HG2	6:S4:255:ARG:HH22	3.20	0.40
7:S5:48:PHE:CG	7:S5:67:PRO:HB3	2.55	0.40
8:S6:122:GLU:O	8:S6:126:ASP:HB3	2.20	0.40
8:S6:34:GLN:HB3	8:S6:52:ILE:HD12	6.13	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:S7:110:GLN:HG2	1:6:811:A:C5	340.22	0.40
9:S7:120:ALA:O	9:S7:124:LYS:HG2	3.04	0.40
9:S7:91:ILE:HD12	9:S7:91:ILE:HA	1.88	0.40
9:S7:62:VAL:O	9:S7:94:ALA:HA	3.04	0.40
1:2:331:A:H5'	10:S8:33:PRO:HA	2.02	0.40
36:1:1093:A:N3	36:1:1096:U:N3	2.68	0.40
36:1:1317:A:C4	36:1:1319:G:C8	3.09	0.40
36:1:164:A:C6	36:1:258:G:C6	3.10	0.40
36:1:2282:U:O2	36:1:2310:U:H4'	2.22	0.40
36:1:2618:G:OP1	47:M0:116:ARG:NE	2.54	0.40
36:1:3054:U:H3	36:1:3088:G:H1	1.70	0.40
36:1:3346:U:H2'	36:1:3347:A:C8	2.56	0.40
36:1:662:U:H2'	36:1:663:C:C6	2.56	0.40
1:2:1038:U:H4'	24:D2:20:THR:HG21	2.03	0.40
1:2:1475:A:H2'	1:2:1476:C:O4'	2.21	0.40
1:2:1483:A:P	1:2:1521:G:H21	2.45	0.40
1:2:32:U:OP1	25:D3:139:LYS:NZ	2.44	0.40
1:2:612:U:H5	1:2:613:G:HO2'	1.67	0.40
1:2:793:A:H5''	1:2:794:U:H6	1.85	0.40
37:3:76:A:H2	56:N0:50:LYS:HZ1	1.67	0.40
36:5:1632:A:H2'	36:5:1633:C:C6	2.56	0.40
70:O4:70:LYS:HD2	36:5:1803:C:O2'	169.23	0.40
73:O7:9:GLY:HA2	36:5:1844:C:O2	148.59	0.40
62:N6:103:LYS:HE3	36:5:221:A:N6	80.48	0.40
36:5:2393:G:O2'	36:5:2982:A:N6	2.46	0.40
66:O0:53:LYS:HZ1	36:5:2552:C:H5	244.08	0.40
78:Q2:39:GLY:HA3	36:5:2765:C:O3'	173.22	0.40
36:5:285:A:O2'	36:5:286:U:H5	2.05	0.40
36:5:68:C:P	36:5:301:G:H21	2.44	0.40
36:5:1895:A:O2'	36:5:3053:G:H4'	2.21	0.40
36:5:3337:G:H2'	36:5:3338:C:O4'	2.21	0.40
36:5:511:G:H2'	36:5:512:U:O4'	2.22	0.40
1:6:1166:A:H2'	1:6:1167:G:O4'	2.21	0.40
8:S6:179:VAL:HG21	1:6:140:A:H1'	327.75	0.40
1:6:1671:A:H2'	1:6:1672:G:O4'	2.21	0.40
1:6:946:U:C4	1:6:947:U:C4	3.10	0.40
1:6:968:U:H2'	1:6:969:C:O4'	2.20	0.40
42:L5:33:ARG:HD2	37:7:7:G:OP1	270.98	0.40
12:C0:8:ARG:HD3	12:C0:79:TYR:OH	5.01	0.40
13:C1:35:TYR:CD1	13:C1:49:ILE:HG12	3.50	0.40
15:C3:84:ILE:HD13	15:C3:89:TYR:HB2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:C5:111:MET:HG2	20:C8:119:ILE:HG23	2.04	0.40
22:D0:51:VAL:HG22	22:D0:94:GLU:H	3.31	0.40
26:D4:99:LYS:O	26:D4:101:GLU:N	2.55	0.40
27:D5:71:ILE:HG13	27:D5:75:LEU:HB2	4.28	0.40
27:D5:89:ILE:HA	27:D5:102:THR:O	2.20	0.40
28:D6:71:LEU:HD22	28:D6:71:LEU:HA	1.89	0.40
30:D8:66:LEU:HA	30:D8:66:LEU:HD23	1.85	0.40
32:E0:14:VAL:O	32:E0:18:THR:HG23	2.21	0.40
39:L2:202:VAL:HG13	39:L2:217:GLN:HG2	2.03	0.40
41:L4:24:ALA:HB2	41:L4:264:SER:HB2	2.03	0.40
41:L4:64:SER:OG	41:L4:73:ARG:O	2.47	0.40
42:L5:64:ILE:HD13	42:L5:105:ILE:HD12	2.03	0.40
42:L5:226:TYR:O	42:L5:230:ASP:N	3.73	0.40
47:M0:52:LEU:HA	47:M0:52:LEU:HD23	1.79	0.40
51:M5:24:ARG:HG2	51:M5:24:ARG:HH11	3.39	0.40
51:M5:84:PRO:O	51:M5:87:GLN:HB2	2.21	0.40
52:M6:50:ASN:HA	52:M6:53:LYS:HG3	3.07	0.40
54:M8:13:SER:C	54:M8:15:HIS:H	2.55	0.40
57:N1:50:LYS:HB3	57:N1:92:ARG:NH1	2.36	0.40
58:N2:37:LEU:HD23	58:N2:41:ILE:HD11	2.03	0.40
59:N3:35:TYR:CE2	59:N3:37:ILE:HG22	2.94	0.40
65:N9:22:LYS:O	65:N9:23:LYS:HB2	4.55	0.40
65:N9:9:ALA:O	65:N9:12:GLN:HG2	2.22	0.40
71:O5:119:LYS:HA	71:O5:119:LYS:HZ3	4.55	0.40
71:O5:88:LEU:HA	71:O5:88:LEU:HD23	1.75	0.40
78:Q2:8:ARG:HG3	78:Q2:9:LYS:N	2.54	0.40
3:S1:144:ARG:NH2	3:S1:207:LEU:O	2.55	0.40
3:S1:69:CYS:HB3	16:C4:114:ARG:HH11	3.44	0.40
3:S1:70:LEU:HB3	3:S1:79:HIS:HB3	5.23	0.40
3:S1:61:LEU:HD11	3:S1:96:LEU:HD13	8.34	0.40
4:S2:143:TYR:CD2	4:S2:147:ASN:HA	3.66	0.40
4:S2:41:LEU:HD21	4:S2:56:ILE:HG22	2.03	0.40
4:S2:61:LEU:HA	4:S2:62:PRO:HD2	1.88	0.40
8:S6:31:ARG:HD2	8:S6:34:GLN:NE2	2.36	0.40
34:SR:248:ASN:HD21	34:SR:249:ARG:NH1	2.20	0.40
36:1:1063:G:C6	36:1:1097:G:C5	3.10	0.40
36:1:953:G:H2'	36:1:1117:G:H5''	2.04	0.40
36:1:1240:A:H61	36:1:1244:A:H5''	1.87	0.40
36:1:1269:U:H2'	36:1:1269:U:O2	2.21	0.40
36:1:1282:G:C6	36:1:1283:C:C4	3.09	0.40
36:1:1895:A:O2'	36:1:3053:G:H4'	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:191:U:H2'	36:1:192:C:H6	1.87	0.40
36:1:2392:C:H5''	36:1:2393:G:OP2	2.20	0.40
36:1:3198:U:O2	46:L9:21:LYS:N	2.55	0.40
36:1:536:U:N3	36:1:556:U:C4	2.89	0.40
36:1:793:C:H5'	64:N8:5:PHE:HB2	2.04	0.40
36:1:992:A:N6	36:1:993:G:O6	2.55	0.40
1:2:1241:G:H2'	1:2:1242:A:C8	2.55	0.40
1:2:412:A:H2	1:2:421:A:H61	1.69	0.40
1:2:463:U:H2'	1:2:464:A:C8	2.56	0.40
36:5:1152:G:N2	36:5:1200:A:C6	2.86	0.40
36:5:1573:G:C5	36:5:1574:C:H1'	2.57	0.40
36:5:1804:A:H2'	36:5:1805:C:O4'	2.21	0.40
36:5:2533:G:OP2	36:5:2533:G:H8	2.04	0.40
36:5:2373:A:N7	36:5:2867:C:H1'	2.36	0.40
69:O3:65:ARG:NH2	36:5:432:G:OP1	203.37	0.40
41:L4:232:SER:O	36:5:694:C:H4'	100.74	0.40
1:6:1263:G:C2	1:6:1264:G:H1'	2.56	0.40
21:C9:37:VAL:HG13	1:6:1503:A:O2'	385.55	0.40
1:6:1515:A:H4'	1:6:1517:U:H5	1.85	0.40
1:6:1619:C:H2'	1:6:1620:C:C6	2.56	0.40
1:6:181:A:H2'	1:6:182:A:O4'	2.21	0.40
25:D3:112:LYS:NZ	1:6:18:C:O3'	346.91	0.40
1:6:397:A:H2'	1:6:398:G:O4'	2.21	0.40
1:6:555:A:C8	1:6:555:A:H3'	2.56	0.40
1:6:901:G:C6	1:6:902:G:C6	3.09	0.40
15:C3:15:ALA:HB2	29:D7:20:LYS:CG	3.29	0.40
17:C5:34:VAL:O	17:C5:42:ARG:HG2	2.25	0.40
17:C5:43:ARG:O	17:C5:47:ARG:HG3	2.20	0.40
20:C8:142:GLY:O	20:C8:145:ARG:HD2	3.20	0.40
20:C8:42:TYR:O	20:C8:46:VAL:HG23	2.49	0.40
21:C9:14:PHE:HZ	21:C9:132:LEU:HG	1.86	0.40
21:C9:53:TRP:HA	21:C9:56:LYS:HB2	2.81	0.40
23:D1:3:ASN:OD1	23:D1:9:VAL:HB	2.21	0.40
24:D2:36:LYS:HD3	24:D2:39:GLN:OE1	2.63	0.40
27:D5:55:PRO:O	27:D5:103:ARG:NH2	8.42	0.40
28:D6:28:LYS:O	28:D6:29:SER:HB3	4.67	0.40
39:L2:136:ILE:HA	39:L2:148:VAL:HG12	2.03	0.40
39:L2:190:ARG:NH1	39:L2:191:LEU:HD11	2.37	0.40
36:1:2163:C:O2'	39:L2:8:GLN:O	2.39	0.40
40:L3:3:HIS:CG	40:L3:3:HIS:O	2.82	0.40
40:L3:84:VAL:HG23	40:L3:163:HIS:C	3.66	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:L4:193:LYS:HE3	41:L4:193:LYS:HB3	2.22	0.40
41:L4:329:PRO:C	41:L4:331:ALA:N	3.01	0.40
41:L4:49:ALA:HA	41:L4:109:TRP:CZ2	2.56	0.40
42:L5:208:MET:HG3	42:L5:223:PHE:CZ	2.56	0.40
36:1:2525:G:H4'	45:L8:49:TYR:OH	2.21	0.40
46:L9:45:PHE:CE1	46:L9:55:VAL:HG13	2.69	0.40
49:M3:122:LYS:HB3	49:M3:122:LYS:NZ	2.36	0.40
50:M4:127:LYS:O	50:M4:131:VAL:HG23	2.94	0.40
50:M4:135:LEU:C	50:M4:137:LYS:H	2.44	0.40
53:M7:95:LEU:HD23	53:M7:95:LEU:HA	2.12	0.40
54:M8:133:LYS:HB2	54:M8:135:GLN:NE2	2.36	0.40
54:M8:98:LYS:HB3	54:M8:99:THR:H	1.70	0.40
55:M9:115:ILE:HD12	55:M9:142:ILE:HD13	2.02	0.40
50:M4:38:ILE:HD13	56:N0:150:PHE:CE2	3.78	0.40
56:N0:31:ALA:HB3	56:N0:37:ALA:HB2	2.03	0.40
60:N4:23:ARG:HE	60:N4:23:ARG:HB3	2.05	0.40
64:N8:42:ARG:HG3	64:N8:43:ILE:N	2.51	0.40
36:1:1327:C:O2'	69:O3:76:GLY:HA2	2.21	0.40
70:O4:104:VAL:HA	70:O4:107:GLU:CD	2.42	0.40
36:1:135:C:H1'	71:O5:93:THR:HB	2.04	0.40
2:S0:169:SER:O	2:S0:173:ILE:HG12	2.20	0.40
4:S2:72:LEU:HD12	4:S2:72:LEU:HA	2.37	0.40
5:S3:80:ALA:HA	5:S3:81:PRO:HD3	1.94	0.40
6:S4:49:ARG:O	6:S4:53:LYS:HD2	4.05	0.40
6:S4:62:LYS:HD2	6:S4:66:MET:HG2	4.78	0.40
7:S5:61:TYR:HE2	7:S5:164:PRO:HG2	4.03	0.40
7:S5:70:VAL:O	7:S5:72:HIS:N	2.56	0.40
9:S7:68:ALA:O	9:S7:72:LYS:HG3	2.53	0.40
10:S8:110:ARG:NH1	10:S8:160:PHE:HB3	3.50	0.40
11:S9:34:PHE:CE1	11:S9:105:LEU:HB3	2.55	0.40
11:S9:134:ILE:HA	11:S9:158:PHE:HA	2.03	0.40
35:SM:102:THR:HG23	35:SM:105:LYS:H	1.86	0.40
35:SM:64:LYS:HD2	35:SM:64:LYS:H	1.85	0.40
34:SR:44:SER:O	34:SR:58:VAL:HG22	2.21	0.40
36:1:1193:A:H2'	36:1:1194:G:O4'	2.22	0.40
36:1:1549:U:H2'	36:1:1550:C:H6	1.87	0.40
36:1:2247:G:H2'	36:1:2248:C:O4'	2.21	0.40
36:1:2259:A:H2'	36:1:2260:U:O4'	2.21	0.40
36:1:2315:G:C2	36:1:2316:G:N7	2.89	0.40
36:1:2510:U:O2'	36:1:2511:A:H5''	2.22	0.40
36:1:2735:U:H2'	36:1:2736:A:H8	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:296:A:O2'	36:1:299:G:H4'	2.21	0.40
36:1:3237:U:H2'	36:1:3238:G:C8	2.57	0.40
36:1:437:G:H2'	36:1:438:A:C8	2.56	0.40
36:1:527:A:H2'	36:1:528:U:O4'	2.21	0.40
36:1:513:G:C2	36:1:579:G:C4	3.09	0.40
1:2:100:A:H2'	1:2:101:U:O4'	2.22	0.40
1:2:1244:A:HO2'	1:2:1245:G:P	2.44	0.40
1:2:1583:A:N1	1:2:1611:A:H5''	2.37	0.40
1:2:850:A:C2	1:2:851:U:C2	3.10	0.40
38:4:118:C:H2'	38:4:119:C:C6	2.57	0.40
38:4:121:U:H2'	38:4:122:U:C6	2.57	0.40
36:5:161:G:H8	36:5:161:G:O5'	2.04	0.40
36:5:1636:U:H3	36:5:1710:C:H4'	1.85	0.40
62:N6:19:TYR:CE2	36:5:216:G:H4'	71.76	0.40
36:5:2660:G:OP1	36:5:2750:U:O2'	2.35	0.40
57:N1:54:HIS:NE2	36:5:2724:U:H4'	229.14	0.40
36:5:2777:G:H5'	36:5:2779:A:OP2	2.22	0.40
78:Q2:56:PRO:HA	36:5:2802:A:C8	183.29	0.40
36:5:2843:U:H5''	36:5:2844:C:OP2	2.21	0.40
36:5:2905:U:H2'	36:5:2906:C:C6	2.56	0.40
36:5:3132:C:H2'	36:5:3133:C:C6	2.56	0.40
36:5:3194:C:C2'	36:5:3195:U:H3'	2.50	0.40
36:5:3222:U:O2'	36:5:3223:A:H5'	2.22	0.40
36:5:573:C:H2'	36:5:574:U:C6	2.56	0.40
36:5:701:G:H2'	36:5:702:C:C6	2.56	0.40
1:6:1017:U:H2'	1:6:1018:U:C6	2.57	0.40
1:6:1315:U:P	1:6:1328:G:H1	2.44	0.40
1:6:1363:U:H6	1:6:1363:U:H2'	1.64	0.40
1:6:1529:C:H2'	1:6:1530:C:C6	2.57	0.40
1:6:1584:G:N2	1:6:1610:G:H2'	2.36	0.40
1:6:778:G:C5	1:6:780:A:C8	3.09	0.40
38:8:104:A:C8	38:8:105:A:C8	3.09	0.40
38:8:15:G:C6	38:8:16:G:C6	3.09	0.40
13:C1:107:VAL:HA	13:C1:108:PRO:HD3	1.84	0.40
1:2:325:G:H4'	13:C1:83:THR:HG21	2.04	0.40
15:C3:27:LYS:HB2	15:C3:28:LEU:H	1.55	0.40
15:C3:33:VAL:HG21	15:C3:66:ILE:HG12	2.04	0.40
16:C4:19:ILE:HD11	16:C4:105:LEU:HD21	2.03	0.40
24:D2:38:LEU:HB3	24:D2:50:PHE:CE1	2.56	0.40
25:D3:91:GLY:C	25:D3:93:LEU:H	2.25	0.40
28:D6:73:TYR:HB3	28:D6:78:ALA:HB2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:D7:36:LYS:HB3	29:D7:42:ASN:O	5.42	0.40
40:L3:232:ARG:HG2	40:L3:233:TRP:CD1	2.74	0.40
41:L4:194:TYR:H	41:L4:194:TYR:HD2	1.70	0.40
41:L4:42:VAL:HG12	41:L4:236:LEU:HD11	2.02	0.40
42:L5:237:GLU:O	42:L5:241:THR:HB	3.05	0.40
42:L5:94:ASN:OD1	42:L5:94:ASN:N	2.92	0.40
36:1:3269:U:H3	43:L6:134:ARG:HH21	1.70	0.40
43:L6:139:LYS:O	43:L6:143:LYS:HG3	2.21	0.40
44:L7:138:TYR:CE2	44:L7:233:GLU:HG2	2.56	0.40
47:M0:99:ILE:HD13	47:M0:99:ILE:H	5.40	0.40
51:M5:36:ILE:HD12	51:M5:106:VAL:HG12	2.03	0.40
36:1:269:G:P	51:M5:44:ARG:HH22	2.44	0.40
54:M8:86:THR:HB	54:M8:105:ARG:HB2	3.19	0.40
57:N1:103:GLN:HE21	57:N1:103:GLN:HB3	1.76	0.40
59:N3:81:GLN:O	59:N3:82:ALA:HB3	2.22	0.40
64:N8:21:ARG:O	64:N8:24:LYS:NZ	2.53	0.40
63:N7:4:PHE:CZ	66:O0:35:ARG:HA	2.57	0.40
66:O0:53:LYS:NZ	36:5:2552:C:H5	243.28	0.40
36:1:361:A:H4'	73:O7:45:ARG:NH2	2.36	0.40
78:Q2:12:CYS:O	78:Q2:18:ARG:HA	2.21	0.40
2:S0:32:HIS:ND1	2:S0:33:GLN:HG3	2.36	0.40
3:S1:87:ARG:NH2	3:S1:89:ASP:OD1	4.66	0.40
4:S2:121:VAL:HB	35:SM:120:GLU:OE1	2.21	0.40
4:S2:56:ILE:CG2	4:S2:61:LEU:HB2	2.91	0.40
6:S4:136:VAL:HG12	6:S4:137:PRO:O	2.94	0.40
7:S5:149:VAL:HG13	7:S5:151:GLY:N	5.28	0.40
7:S5:65:ARG:HA	7:S5:66:GLN:HA	1.85	0.40
8:S6:51:LYS:HB3	8:S6:112:VAL:HG13	4.62	0.40
8:S6:73:ILE:O	8:S6:97:VAL:HG12	2.20	0.40
8:S6:77:LEU:HD12	8:S6:95:LYS:HB2	2.83	0.40
9:S7:141:ARG:HG2	24:D2:51:GLU:CD	3.24	0.40
11:S9:107:ARG:O	11:S9:147:MET:HA	2.21	0.40
11:S9:154:LYS:HB2	11:S9:154:LYS:HE3	1.86	0.40
36:1:1026:A:H2'	36:1:1027:A:C8	2.57	0.40
36:1:1083:G:H2'	36:1:1084:A:C8	2.56	0.40
36:1:1083:G:C2	36:1:1084:A:C4	3.10	0.40
36:1:1171:G:H2'	36:1:1172:G:O4'	2.22	0.40
36:1:1238:C:H1'	36:1:1251:A:H61	1.86	0.40
36:1:1764:U:H3'	36:1:1765:U:C4'	2.51	0.40
36:1:214:G:H4'	62:N6:10:SER:O	2.21	0.40
36:1:2366:C:H2'	36:1:2367:A:C8	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:2630:C:C5	57:N1:4:SER:HB2	2.56	0.40
36:1:3144:G:N2	36:1:3145:C:H1'	2.36	0.40
36:1:430:U:H2'	36:1:431:U:O4'	2.22	0.40
36:1:439:C:O2	36:1:439:C:H2'	2.21	0.40
36:1:600:G:N2	36:1:604:G:C6	2.90	0.40
36:1:980:A:H2'	36:1:981:U:C1'	2.52	0.40
1:2:1228:G:OP2	14:C2:119:SER:OG	2.39	0.40
1:2:1244:A:O2'	1:2:1245:G:OP1	2.37	0.40
1:2:1504:G:C6	1:2:1505:A:C6	3.09	0.40
1:2:1615:C:C5	7:S5:81:ARG:HA	2.57	0.40
1:2:325:G:H2'	1:2:326:G:C8	2.57	0.40
37:3:36:C:O2'	37:3:37:G:H5'	2.22	0.40
67:O1:26:LYS:NZ	36:5:1456:A:N7	166.89	0.40
36:5:158:G:N2	36:5:264:G:H1'	2.36	0.40
36:5:1658:G:H2'	36:5:1659:U:H6	1.86	0.40
36:5:3028:G:H2'	36:5:3029:A:C8	2.56	0.40
36:5:3087:A:H4'	36:5:3375:A:H61	1.87	0.40
36:5:3234:A:N6	36:5:3235:C:C4	2.90	0.40
36:5:540:U:H2'	36:5:541:U:O4'	2.21	0.40
36:5:361:A:N3	36:5:814:U:H1'	2.37	0.40
36:5:822:G:H2'	36:5:823:C:O4'	2.21	0.40
1:6:1117:U:H2'	1:6:1118:G:H8	1.87	0.40
1:6:717:C:O2'	1:6:718:U:OP1	2.39	0.40
14:C2:78:LEU:HA	14:C2:78:LEU:HD23	1.92	0.40
16:C4:48:VAL:HG22	16:C4:49:LYS:H	3.08	0.40
17:C5:21:ASP:O	17:C5:23:GLU:N	2.54	0.40
17:C5:84:ILE:H	17:C5:84:ILE:HD12	4.55	0.40
20:C8:56:LYS:HG2	20:C8:60:GLU:CD	2.41	0.40
22:D0:96:PRO:HB2	22:D0:97:VAL:H	1.65	0.40
24:D2:7:LEU:HD11	24:D2:37:PHE:HD2	3.68	0.40
25:D3:87:VAL:HG22	25:D3:124:VAL:HG21	2.03	0.40
26:D4:117:LYS:HG2	1:6:159:U:H5'	332.85	0.40
23:D1:64:GLU:CG	29:D7:3:LEU:HG	3.42	0.40
40:L3:106:TRP:CG	40:L3:130:PHE:CE1	3.32	0.40
40:L3:339:ARG:NH1	40:L3:342:LEU:HD21	2.36	0.40
40:L3:80:ASP:OD2	40:L3:314:TYR:OH	2.38	0.40
41:L4:316:ASN:HA	41:L4:317:PRO:HD3	2.14	0.40
41:L4:68:GLY:O	36:5:2401:A:H4'	171.06	0.40
45:L8:153:ILE:HD13	45:L8:167:PRO:HD3	2.02	0.40
45:L8:75:ILE:HG22	45:L8:76:ALA:N	2.37	0.40
46:L9:189:GLU:O	46:L9:191:LEU:N	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M5:80:THR:HG21	51:M5:87:GLN:HA	2.03	0.40
56:N0:155:ARG:HB2	56:N0:172:TYR:HB2	2.44	0.40
56:N0:71:LYS:O	56:N0:73:LYS:HG3	2.22	0.40
59:N3:27:ASP:HA	59:N3:113:ALA:O	2.57	0.40
60:N4:14:TYR:O	60:N4:17:ARG:HB3	2.21	0.40
61:N5:77:GLU:HB3	61:N5:78:ASP:OD1	2.21	0.40
63:N7:110:ALA:O	63:N7:113:VAL:HB	2.94	0.40
63:N7:126:LYS:O	63:N7:128:GLN:N	2.45	0.40
66:O0:41:LEU:HA	66:O0:41:LEU:HD22	1.94	0.40
68:O2:118:LYS:HG2	68:O2:119:VAL:N	3.03	0.40
69:O3:67:MET:HE1	69:O3:87:ASN:HB2	5.41	0.40
36:1:1694:U:O2'	70:O4:25:THR:O	2.34	0.40
36:1:1597:C:OP1	70:O4:31:ARG:HD2	2.22	0.40
77:Q1:8:LYS:HD3	77:Q1:12:ARG:HH21	1.85	0.40
78:Q2:74:CYS:O	78:Q2:78:LYS:N	2.55	0.40
3:S1:86:LEU:HB3	3:S1:98:THR:OG1	2.22	0.40
6:S4:31:PRO:HG2	6:S4:38:LEU:O	2.97	0.40
7:S5:133:VAL:HG22	7:S5:198:LEU:HD13	2.05	0.40
9:S7:89:HIS:ND1	9:S7:168:SER:OG	3.45	0.40
11:S9:16:LYS:HB3	11:S9:16:LYS:HE3	4.92	0.40
34:SR:111:MET:N	34:SR:125:GLY:O	2.93	0.40
34:SR:22:SER:HB3	34:SR:36:ALA:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:1:3153:U:O4	34:sR:79:TYR:OH[2_656]	2.12	0.08

## 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	S0	204/251 (81%)	151 (74%)	30 (15%)	23 (11%)	0	3
2	s0	204/251 (81%)	164 (80%)	25 (12%)	15 (7%)	1	10
3	S1	212/254 (84%)	152 (72%)	38 (18%)	22 (10%)	1	4
3	s1	214/254 (84%)	168 (78%)	33 (15%)	13 (6%)	2	14
4	S2	215/253 (85%)	168 (78%)	33 (15%)	14 (6%)	1	13
4	s2	215/253 (85%)	168 (78%)	33 (15%)	14 (6%)	1	13
5	S3	221/239 (92%)	187 (85%)	26 (12%)	8 (4%)	4	28
5	s3	221/239 (92%)	176 (80%)	35 (16%)	10 (4%)	3	21
6	S4	258/260 (99%)	202 (78%)	46 (18%)	10 (4%)	4	25
6	s4	258/260 (99%)	189 (73%)	53 (20%)	16 (6%)	2	14
7	S5	204/224 (91%)	155 (76%)	37 (18%)	12 (6%)	2	15
7	s5	204/224 (91%)	160 (78%)	29 (14%)	15 (7%)	1	10
8	S6	224/236 (95%)	193 (86%)	22 (10%)	9 (4%)	4	24
8	s6	216/236 (92%)	182 (84%)	22 (10%)	12 (6%)	2	16
9	S7	182/189 (96%)	134 (74%)	27 (15%)	21 (12%)	0	3
9	s7	184/189 (97%)	152 (83%)	20 (11%)	12 (6%)	1	13
10	S8	184/200 (92%)	151 (82%)	24 (13%)	9 (5%)	3	19
10	s8	184/200 (92%)	157 (85%)	20 (11%)	7 (4%)	4	26
11	S9	183/196 (93%)	149 (81%)	23 (13%)	11 (6%)	2	15
11	s9	183/196 (93%)	147 (80%)	27 (15%)	9 (5%)	3	19
12	C0	82/96 (85%)	71 (87%)	8 (10%)	3 (4%)	4	27
12	c0	82/96 (85%)	56 (68%)	16 (20%)	10 (12%)	0	2
13	C1	143/155 (92%)	116 (81%)	18 (13%)	9 (6%)	2	13
13	c1	143/155 (92%)	115 (80%)	21 (15%)	7 (5%)	3	19
14	C2	107/142 (75%)	70 (65%)	27 (25%)	10 (9%)	1	6
14	c2	107/142 (75%)	75 (70%)	22 (21%)	10 (9%)	1	6
15	C3	148/150 (99%)	124 (84%)	20 (14%)	4 (3%)	6	35
15	c3	148/150 (99%)	116 (78%)	23 (16%)	9 (6%)	2	14
16	C4	125/136 (92%)	81 (65%)	33 (26%)	11 (9%)	1	7
16	c4	126/136 (93%)	96 (76%)	15 (12%)	15 (12%)	0	3
17	C5	121/141 (86%)	95 (78%)	12 (10%)	14 (12%)	0	3
17	c5	128/141 (91%)	98 (77%)	19 (15%)	11 (9%)	1	7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
18	C6	139/142 (98%)	116 (84%)	14 (10%)	9 (6%)	1	13
18	c6	140/142 (99%)	118 (84%)	13 (9%)	9 (6%)	2	13
19	C7	116/136 (85%)	82 (71%)	23 (20%)	11 (10%)	1	6
19	c7	113/136 (83%)	93 (82%)	13 (12%)	7 (6%)	2	14
20	C8	143/145 (99%)	116 (81%)	17 (12%)	10 (7%)	1	11
20	c8	143/145 (99%)	120 (84%)	16 (11%)	7 (5%)	3	19
21	C9	141/143 (99%)	115 (82%)	18 (13%)	8 (6%)	2	16
21	c9	141/143 (99%)	120 (85%)	15 (11%)	6 (4%)	3	23
22	D0	105/120 (88%)	89 (85%)	9 (9%)	7 (7%)	1	12
22	d0	108/120 (90%)	89 (82%)	14 (13%)	5 (5%)	3	21
23	D1	85/87 (98%)	66 (78%)	11 (13%)	8 (9%)	1	6
23	d1	85/87 (98%)	63 (74%)	17 (20%)	5 (6%)	2	15
24	D2	127/129 (98%)	116 (91%)	8 (6%)	3 (2%)	7	38
24	d2	127/129 (98%)	107 (84%)	17 (13%)	3 (2%)	7	38
25	D3	142/144 (99%)	109 (77%)	21 (15%)	12 (8%)	1	7
25	d3	142/144 (99%)	117 (82%)	20 (14%)	5 (4%)	4	29
26	D4	132/134 (98%)	109 (83%)	13 (10%)	10 (8%)	1	9
26	d4	132/134 (98%)	110 (83%)	15 (11%)	7 (5%)	2	17
27	D5	68/107 (64%)	48 (71%)	13 (19%)	7 (10%)	1	4
27	d5	67/107 (63%)	51 (76%)	15 (22%)	1 (2%)	13	49
28	D6	95/97 (98%)	56 (59%)	21 (22%)	18 (19%)	0	1
28	d6	95/97 (98%)	66 (70%)	20 (21%)	9 (10%)	1	6
29	D7	79/81 (98%)	61 (77%)	15 (19%)	3 (4%)	4	26
29	d7	79/81 (98%)	61 (77%)	14 (18%)	4 (5%)	2	19
30	D8	61/66 (92%)	50 (82%)	9 (15%)	2 (3%)	5	30
30	d8	61/66 (92%)	47 (77%)	11 (18%)	3 (5%)	3	19
31	D9	51/55 (93%)	39 (76%)	9 (18%)	3 (6%)	2	15
31	d9	51/55 (93%)	41 (80%)	8 (16%)	2 (4%)	4	25
32	E0	58/62 (94%)	46 (79%)	10 (17%)	2 (3%)	5	29
32	e0	60/62 (97%)	40 (67%)	12 (20%)	8 (13%)	0	2
33	E1	69/76 (91%)	35 (51%)	17 (25%)	17 (25%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
33	e1	74/76 (97%)	36 (49%)	21 (28%)	17 (23%)	0	0
34	SR	316/318 (99%)	263 (83%)	36 (11%)	17 (5%)	2	17
34	sR	316/318 (99%)	252 (80%)	53 (17%)	11 (4%)	4	29
35	SM	120/273 (44%)	91 (76%)	18 (15%)	11 (9%)	1	6
35	sM	61/273 (22%)	42 (69%)	11 (18%)	8 (13%)	0	2
39	L2	250/253 (99%)	213 (85%)	27 (11%)	10 (4%)	4	24
39	l2	250/253 (99%)	205 (82%)	30 (12%)	15 (6%)	2	15
40	L3	384/386 (100%)	316 (82%)	49 (13%)	19 (5%)	3	19
40	l3	384/386 (100%)	330 (86%)	39 (10%)	15 (4%)	4	25
41	L4	359/361 (99%)	291 (81%)	46 (13%)	22 (6%)	2	14
41	l4	359/361 (99%)	296 (82%)	43 (12%)	20 (6%)	2	16
42	L5	294/296 (99%)	232 (79%)	44 (15%)	18 (6%)	2	14
42	l5	292/296 (99%)	248 (85%)	38 (13%)	6 (2%)	9	42
43	L6	152/175 (87%)	122 (80%)	27 (18%)	3 (2%)	9	43
43	l6	153/175 (87%)	125 (82%)	24 (16%)	4 (3%)	7	36
44	L7	220/243 (90%)	185 (84%)	24 (11%)	11 (5%)	3	19
44	l7	221/243 (91%)	192 (87%)	21 (10%)	8 (4%)	4	28
45	L8	231/255 (91%)	190 (82%)	30 (13%)	11 (5%)	3	20
45	l8	229/255 (90%)	174 (76%)	40 (18%)	15 (7%)	1	12
46	L9	189/191 (99%)	145 (77%)	29 (15%)	15 (8%)	1	8
46	l9	189/191 (99%)	158 (84%)	24 (13%)	7 (4%)	4	27
47	M0	207/220 (94%)	161 (78%)	37 (18%)	9 (4%)	3	23
47	m0	209/220 (95%)	169 (81%)	29 (14%)	11 (5%)	2	17
48	M1	167/173 (96%)	128 (77%)	27 (16%)	12 (7%)	1	11
48	m1	167/173 (96%)	145 (87%)	9 (5%)	13 (8%)	1	9
49	M3	191/198 (96%)	153 (80%)	30 (16%)	8 (4%)	3	23
49	m3	192/198 (97%)	152 (79%)	27 (14%)	13 (7%)	1	12
50	M4	134/137 (98%)	116 (87%)	12 (9%)	6 (4%)	3	21
50	m4	135/137 (98%)	117 (87%)	15 (11%)	3 (2%)	8	41
51	M5	201/203 (99%)	175 (87%)	20 (10%)	6 (3%)	5	33
51	m5	201/203 (99%)	178 (89%)	18 (9%)	5 (2%)	7	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
52	M6	195/198 (98%)	174 (89%)	12 (6%)	9 (5%)	3	21
52	m6	195/198 (98%)	172 (88%)	20 (10%)	3 (2%)	13	49
53	M7	181/183 (99%)	151 (83%)	25 (14%)	5 (3%)	6	34
53	m7	153/183 (84%)	137 (90%)	14 (9%)	2 (1%)	15	52
54	M8	183/185 (99%)	156 (85%)	20 (11%)	7 (4%)	4	26
54	m8	183/185 (99%)	154 (84%)	21 (12%)	8 (4%)	3	22
55	M9	186/188 (99%)	164 (88%)	17 (9%)	5 (3%)	6	35
55	m9	186/188 (99%)	164 (88%)	20 (11%)	2 (1%)	17	57
56	N0	170/172 (99%)	151 (89%)	15 (9%)	4 (2%)	7	38
56	n0	170/172 (99%)	159 (94%)	9 (5%)	2 (1%)	16	54
57	N1	157/159 (99%)	131 (83%)	20 (13%)	6 (4%)	4	26
57	n1	157/159 (99%)	138 (88%)	16 (10%)	3 (2%)	10	45
58	N2	98/120 (82%)	73 (74%)	20 (20%)	5 (5%)	2	19
58	n2	96/120 (80%)	80 (83%)	13 (14%)	3 (3%)	5	32
59	N3	134/136 (98%)	117 (87%)	12 (9%)	5 (4%)	4	27
59	n3	134/136 (98%)	122 (91%)	4 (3%)	8 (6%)	2	15
60	N4	96/155 (62%)	66 (69%)	21 (22%)	9 (9%)	1	6
60	n4	133/155 (86%)	102 (77%)	23 (17%)	8 (6%)	2	15
61	N5	119/141 (84%)	103 (87%)	12 (10%)	4 (3%)	5	29
61	n5	118/141 (84%)	91 (77%)	17 (14%)	10 (8%)	1	7
62	N6	124/126 (98%)	104 (84%)	19 (15%)	1 (1%)	24	62
62	n6	124/126 (98%)	102 (82%)	17 (14%)	5 (4%)	4	24
63	N7	133/135 (98%)	112 (84%)	12 (9%)	9 (7%)	1	12
63	n7	133/135 (98%)	100 (75%)	26 (20%)	7 (5%)	2	17
64	N8	146/148 (99%)	113 (77%)	26 (18%)	7 (5%)	3	20
64	n8	146/148 (99%)	110 (75%)	25 (17%)	11 (8%)	1	10
65	N9	56/58 (97%)	46 (82%)	8 (14%)	2 (4%)	4	28
65	n9	56/58 (97%)	41 (73%)	9 (16%)	6 (11%)	0	4
66	O0	95/104 (91%)	83 (87%)	11 (12%)	1 (1%)	17	57
66	o0	98/104 (94%)	86 (88%)	10 (10%)	2 (2%)	9	43
67	O1	107/112 (96%)	96 (90%)	6 (6%)	5 (5%)	3	20

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
67	o1	107/112 (96%)	85 (79%)	15 (14%)	7 (6%)	1	13
68	O2	125/129 (97%)	110 (88%)	11 (9%)	4 (3%)	5	31
68	o2	125/129 (97%)	105 (84%)	16 (13%)	4 (3%)	5	31
69	O3	104/106 (98%)	96 (92%)	6 (6%)	2 (2%)	10	45
69	o3	104/106 (98%)	93 (89%)	9 (9%)	2 (2%)	10	45
70	O4	110/120 (92%)	96 (87%)	11 (10%)	3 (3%)	6	35
70	o4	110/120 (92%)	88 (80%)	18 (16%)	4 (4%)	4	28
71	O5	117/119 (98%)	95 (81%)	12 (10%)	10 (8%)	1	7
71	o5	117/119 (98%)	94 (80%)	21 (18%)	2 (2%)	11	47
72	O6	97/99 (98%)	76 (78%)	12 (12%)	9 (9%)	1	6
72	o6	97/99 (98%)	76 (78%)	11 (11%)	10 (10%)	1	4
73	O7	85/87 (98%)	69 (81%)	14 (16%)	2 (2%)	7	38
73	o7	85/87 (98%)	70 (82%)	14 (16%)	1 (1%)	16	54
74	O8	75/77 (97%)	66 (88%)	7 (9%)	2 (3%)	6	35
74	o8	75/77 (97%)	61 (81%)	9 (12%)	5 (7%)	1	12
75	O9	48/50 (96%)	35 (73%)	10 (21%)	3 (6%)	2	13
75	o9	48/50 (96%)	39 (81%)	9 (19%)	0	100	100
76	Q0	50/52 (96%)	43 (86%)	6 (12%)	1 (2%)	9	43
76	q0	50/52 (96%)	48 (96%)	1 (2%)	1 (2%)	9	43
77	Q1	23/25 (92%)	21 (91%)	2 (9%)	0	100	100
77	q1	23/25 (92%)	20 (87%)	2 (9%)	1 (4%)	3	23
78	Q2	103/105 (98%)	85 (82%)	13 (13%)	5 (5%)	3	19
78	q2	103/105 (98%)	88 (85%)	9 (9%)	6 (6%)	2	15
79	Q3	89/91 (98%)	70 (79%)	13 (15%)	6 (7%)	1	12
79	q3	89/91 (98%)	79 (89%)	9 (10%)	1 (1%)	17	57
81	p0	117/311 (38%)	97 (83%)	14 (12%)	6 (5%)	2	19
83	f	146/157 (93%)	91 (62%)	35 (24%)	20 (14%)	0	2
All	All	22316/24284 (92%)	18108 (81%)	2992 (13%)	1216 (5%)	2	17

All (1216) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	S0	95	ALA

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Mol	Chain	Res	Type
2	S0	158	VAL
2	S0	190	ASP
2	S0	192	THR
3	S1	35	PRO
3	S1	37	THR
3	S1	132	ASP
3	S1	206	PRO
3	S1	213	ARG
4	S2	48	GLY
4	S2	106	ASP
5	S3	217	ILE
5	S3	220	PRO
6	S4	12	LEU
7	S5	63	GLN
7	S5	98	MET
8	S6	122	GLU
8	S6	173	PRO
9	S7	31	SER
9	S7	32	PRO
9	S7	64	VAL
9	S7	116	ARG
9	S7	186	PRO
11	S9	93	LEU
11	S9	98	ALA
11	S9	134	ILE
13	C1	7	VAL
13	C1	55	ASP
16	C4	42	VAL
17	C5	54	ALA
17	C5	69	GLU
17	C5	80	MET
17	C5	87	PRO
17	C5	125	PRO
18	C6	41	PRO
18	C6	58	ASP
18	C6	97	VAL
18	C6	114	ARG
19	C7	85	VAL
19	C7	86	PRO
19	C7	88	VAL
19	C7	99	VAL
19	C7	124	VAL

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Mol	Chain	Res	Type
20	C8	76	PRO
20	C8	92	ILE
21	C9	11	ALA
21	C9	31	PRO
21	C9	53	TRP
21	C9	67	MET
21	C9	69	LYS
23	D1	4	ASP
24	D2	83	ILE
25	D3	36	THR
25	D3	114	LYS
25	D3	144	ARG
27	D5	44	GLN
27	D5	71	ILE
27	D5	86	GLU
28	D6	45	VAL
28	D6	84	VAL
29	D7	57	GLU
31	D9	11	PRO
32	E0	47	VAL
33	E1	84	VAL
33	E1	98	VAL
33	E1	102	VAL
33	E1	103	LEU
34	SR	238	ASP
34	SR	318	ALA
35	SM	64	LYS
39	L2	144	ASN
40	L3	3	HIS
40	L3	5	LYS
41	L4	4	PRO
41	L4	130	ALA
41	L4	131	VAL
41	L4	220	ARG
41	L4	293	SER
41	L4	341	SER
42	L5	178	ASN
42	L5	223	PHE
42	L5	258	LYS
43	L6	98	VAL
44	L7	24	GLU
44	L7	72	ALA

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
45	L8	25	PRO
45	L8	31	PRO
45	L8	122	LYS
45	L8	226	TYR
46	L9	50	ASN
46	L9	107	ASP
47	M0	6	ALA
47	M0	113	GLN
47	M0	130	ASP
47	M0	219	ALA
48	M1	11	ASP
49	M3	129	ASN
50	M4	10	SER
51	M5	74	PRO
51	M5	75	VAL
52	M6	110	PRO
52	M6	111	PRO
52	M6	149	TYR
53	M7	157	VAL
57	N1	159	PHE
58	N2	51	GLY
59	N3	54	LEU
60	N4	81	PRO
61	N5	48	SER
63	N7	17	ARG
63	N7	30	ASP
63	N7	125	GLY
63	N7	128	GLN
64	N8	66	ALA
64	N8	96	LYS
64	N8	97	GLU
67	O1	21	HIS
72	O6	28	TYR
72	O6	33	ALA
74	O8	33	LYS
75	O9	4	GLN
78	Q2	96	GLU
78	Q2	100	LYS
2	s0	29	VAL
2	s0	66	ALA
2	s0	68	PRO
2	s0	164	ASN

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Mol	Chain	Res	Type
2	s0	189	VAL
2	s0	194	PRO
3	s1	158	SER
3	s1	218	LEU
4	s2	164	SER
5	s3	217	ILE
5	s3	220	PRO
6	s4	53	LYS
6	s4	163	ASP
6	s4	196	VAL
7	s5	43	PHE
7	s5	184	PHE
7	s5	204	GLY
8	s6	152	ASP
8	s6	153	VAL
9	s7	5	GLN
9	s7	74	GLN
9	s7	106	SER
9	s7	131	PHE
12	c0	2	LEU
12	c0	32	HIS
12	c0	83	PRO
15	c3	66	ILE
16	c4	51	ASP
16	c4	91	THR
16	c4	98	GLY
17	c5	80	MET
17	c5	125	PRO
18	c6	39	VAL
18	c6	42	GLU
18	c6	97	VAL
18	c6	112	TYR
18	c6	113	ASP
18	c6	116	LEU
19	c7	67	ARG
19	c7	88	VAL
19	c7	99	VAL
20	c8	8	GLN
21	c9	33	TYR
24	d2	68	ARG
25	d3	47	SER
25	d3	138	GLU

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Mol	Chain	Res	Type
26	d4	33	ALA
26	d4	35	VAL
26	d4	53	ASP
28	d6	28	LYS
28	d6	29	SER
31	d9	6	VAL
32	e0	60	PRO
33	e1	87	THR
33	e1	92	LYS
33	e1	98	VAL
33	e1	107	LYS
33	e1	110	ALA
34	sR	160	GLU
34	sR	165	ASP
34	sR	226	ALA
35	sM	50	ASN
39	l2	24	GLN
39	l2	142	ASP
39	l2	194	ASN
40	l3	139	GLN
40	l3	170	PRO
40	l3	300	ARG
40	l3	347	SER
41	l4	142	VAL
41	l4	174	ALA
41	l4	252	GLU
41	l4	272	VAL
41	l4	301	PRO
41	l4	329	PRO
41	l4	342	LYS
41	l4	361	HIS
42	l5	178	ASN
44	l7	193	PRO
45	l8	25	PRO
45	l8	81	THR
46	l9	62	ARG
46	l9	167	VAL
47	m0	82	ARG
47	m0	145	LYS
48	m1	8	PRO
48	m1	108	GLU
49	m3	17	HIS

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Mol	Chain	Res	Type
49	m3	134	GLU
50	m4	63	VAL
51	m5	76	PRO
52	m6	5	PRO
52	m6	110	PRO
54	m8	95	GLU
54	m8	99	THR
57	n1	135	PRO
58	n2	49	ASN
58	n2	105	LEU
59	n3	42	SER
60	n4	26	SER
60	n4	63	ILE
60	n4	75	THR
60	n4	76	VAL
61	n5	42	ARG
61	n5	92	LYS
63	n7	56	LYS
64	n8	30	GLY
64	n8	47	LYS
64	n8	48	TYR
65	n9	21	ILE
65	n9	23	LYS
65	n9	24	PRO
67	o1	5	LYS
67	o1	45	GLY
67	o1	84	ASP
68	o2	6	HIS
69	o3	60	ARG
69	o3	94	PHE
72	o6	64	SER
72	o6	98	ARG
76	q0	78	ILE
78	q2	32	LYS
78	q2	33	ALA
83	f	52	HIS
83	f	54	HIS
83	f	64	ILE
83	f	109	THR
83	f	111	ASP
2	S0	4	PRO
2	S0	5	ALA

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Mol	Chain	Res	Type
2	S0	39	ASN
2	S0	111	ILE
2	S0	191	ARG
2	S0	202	TYR
3	S1	21	VAL
3	S1	62	LYS
3	S1	63	GLY
3	S1	177	GLN
4	S2	91	ARG
4	S2	107	SER
5	S3	81	PRO
5	S3	216	PRO
6	S4	104	ASP
6	S4	150	PRO
6	S4	195	ILE
7	S5	39	GLU
7	S5	43	PHE
7	S5	204	GLY
8	S6	149	LYS
8	S6	150	GLU
8	S6	174	LYS
9	S7	36	ALA
9	S7	74	GLN
9	S7	129	LEU
10	S8	52	ASN
10	S8	149	SER
11	S9	167	ALA
11	S9	169	PRO
13	C1	4	GLU
13	C1	19	ILE
13	C1	144	ALA
14	C2	69	ALA
14	C2	91	VAL
14	C2	93	ASP
14	C2	119	SER
14	C2	126	TRP
15	C3	22	ALA
16	C4	12	GLN
16	C4	124	ASP
16	C4	125	SER
16	C4	126	THR
17	C5	11	VAL

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Mol	Chain	Res	Type
17	C5	22	LEU
17	C5	48	GLY
19	C7	98	GLY
19	C7	122	ILE
20	C8	14	ILE
21	C9	10	ALA
21	C9	116	ILE
22	D0	16	GLN
22	D0	17	GLN
22	D0	49	ASN
22	D0	96	PRO
22	D0	119	ALA
25	D3	39	LYS
25	D3	112	LYS
25	D3	137	LYS
25	D3	138	GLU
26	D4	33	ALA
26	D4	54	ALA
26	D4	100	VAL
27	D5	41	ILE
28	D6	4	LYS
28	D6	16	GLY
28	D6	18	VAL
28	D6	36	ILE
28	D6	82	ARG
28	D6	86	VAL
29	D7	62	ILE
30	D8	36	THR
31	D9	20	GLN
32	E0	51	ASN
33	E1	87	THR
33	E1	118	ARG
33	E1	127	GLY
33	E1	144	CYS
33	E1	145	HIS
34	SR	51	ASP
34	SR	98	GLU
34	SR	135	THR
34	SR	160	GLU
34	SR	161	LYS
35	SM	52	PRO
35	SM	87	THR

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Mol	Chain	Res	Type
35	SM	89	ARG
39	L2	13	GLY
39	L2	250	GLN
40	L3	4	ARG
40	L3	120	LYS
40	L3	121	ASN
40	L3	142	ALA
40	L3	155	ALA
40	L3	171	LEU
40	L3	174	LYS
40	L3	221	THR
40	L3	347	SER
40	L3	360	ASP
41	L4	15	ALA
41	L4	83	GLY
41	L4	305	ALA
41	L4	311	HIS
42	L5	7	ALA
42	L5	212	ALA
42	L5	221	GLU
42	L5	253	PHE
44	L7	55	TYR
44	L7	80	GLN
44	L7	158	LYS
44	L7	159	GLN
45	L8	39	ALA
46	L9	38	LEU
46	L9	59	ASN
47	M0	78	THR
47	M0	117	GLY
48	M1	8	PRO
48	M1	114	ILE
48	M1	152	HIS
49	M3	47	ALA
49	M3	141	ALA
50	M4	8	LYS
50	M4	36	VAL
51	M5	81	TYR
51	M5	110	ALA
51	M5	184	LYS
52	M6	86	GLY
52	M6	112	TYR

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Mol	Chain	Res	Type
52	M6	113	ASP
53	M7	109	ALA
54	M8	74	GLU
54	M8	98	LYS
54	M8	166	LEU
55	M9	53	LYS
56	N0	24	LEU
56	N0	154	HIS
57	N1	122	GLN
59	N3	46	LEU
61	N5	24	LEU
61	N5	77	GLU
61	N5	117	ASN
63	N7	16	GLY
64	N8	117	ARG
67	O1	6	ASP
67	O1	99	ALA
69	O3	59	VAL
69	O3	60	ARG
70	O4	18	ASN
71	O5	35	LYS
71	O5	71	LYS
71	O5	118	ILE
72	O6	3	VAL
72	O6	78	GLY
72	O6	99	ARG
78	Q2	30	ALA
78	Q2	34	SER
78	Q2	60	LYS
79	Q3	3	LYS
79	Q3	58	SER
2	s0	4	PRO
2	s0	27	ARG
2	s0	65	ALA
3	s1	43	VAL
3	s1	206	PRO
3	s1	224	ASP
4	s2	93	GLY
4	s2	144	TRP
4	s2	149	GLY
4	s2	249	ALA
5	s3	115	ILE

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Mol	Chain	Res	Type
5	s3	211	PRO
6	s4	3	ARG
6	s4	104	ASP
6	s4	118	GLU
6	s4	195	ILE
7	s5	28	PRO
7	s5	81	ARG
7	s5	155	ALA
8	s6	19	ASP
8	s6	58	LYS
9	s7	63	PRO
9	s7	144	VAL
10	s8	52	ASN
10	s8	101	ILE
11	s9	115	LYS
11	s9	121	SER
12	c0	23	ALA
12	c0	82	LEU
13	c1	55	ASP
14	c2	131	ASP
15	c3	43	LYS
15	c3	137	PRO
15	c3	140	LYS
16	c4	14	PHE
17	c5	17	TYR
19	c7	116	LYS
20	c8	14	ILE
20	c8	60	GLU
21	c9	29	GLU
22	d0	15	GLN
22	d0	119	ALA
23	d1	11	LEU
23	d1	85	TYR
24	d2	78	ARG
25	d3	137	LYS
26	d4	52	LYS
26	d4	54	ALA
28	d6	5	ARG
28	d6	13	LYS
28	d6	34	LYS
29	d7	59	CYS
30	d8	33	LEU

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Mol	Chain	Res	Type
32	e0	47	VAL
32	e0	49	LEU
32	e0	50	VAL
33	e1	84	VAL
33	e1	102	VAL
34	sR	163	ASP
35	sM	47	ALA
35	sM	55	SER
39	l2	170	ALA
40	l3	129	ALA
40	l3	187	SER
40	l3	385	LYS
40	l3	386	ASP
41	l4	15	ALA
41	l4	90	PHE
41	l4	154	THR
41	l4	233	LEU
41	l4	268	ALA
41	l4	338	LYS
42	l5	260	PHE
44	l7	206	LYS
45	l8	26	LEU
45	l8	203	VAL
45	l8	237	ILE
46	l9	144	ILE
46	l9	175	PHE
47	m0	117	GLY
47	m0	151	GLY
47	m0	196	PHE
48	m1	94	ARG
48	m1	95	ASN
48	m1	114	ILE
48	m1	115	LYS
48	m1	145	LYS
48	m1	167	TYR
49	m3	13	HIS
49	m3	47	ALA
49	m3	141	ALA
49	m3	193	ALA
50	m4	74	ARG
54	m8	84	VAL
54	m8	98	LYS

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Mol	Chain	Res	Type
54	m8	116	LYS
54	m8	168	THR
55	m9	36	ASN
56	n0	17	GLU
57	n1	136	ARG
59	n3	28	ASN
59	n3	134	GLY
60	n4	83	THR
61	n5	24	LEU
61	n5	44	PRO
61	n5	55	ASN
62	n6	83	ASP
62	n6	84	LYS
63	n7	128	GLN
63	n7	129	TRP
63	n7	130	PHE
64	n8	76	ASP
64	n8	129	PHE
65	n9	5	LYS
66	o0	10	ILE
66	o0	85	PHE
67	o1	91	SER
68	o2	27	ARG
71	o5	43	LYS
71	o5	119	LYS
72	o6	12	ASN
72	o6	20	MET
72	o6	33	ALA
72	o6	35	ASN
72	o6	79	SER
73	o7	65	ARG
74	o8	18	ALA
74	o8	46	ARG
79	q3	51	ALA
81	p0	68	SER
81	p0	93	LEU
83	f	51	LYS
83	f	148	ILE
2	S0	30	GLN
2	S0	66	ALA
2	S0	78	SER
2	S0	161	PRO

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Mol	Chain	Res	Type
3	S1	48	VAL
3	S1	51	SER
3	S1	58	SER
3	S1	81	PHE
3	S1	209	ASN
4	S2	95	ARG
5	S3	62	ASN
5	S3	205	ALA
5	S3	218	LEU
6	S4	205	PHE
6	S4	214	LEU
6	S4	234	PRO
7	S5	26	ALA
7	S5	58	LEU
7	S5	64	VAL
7	S5	97	LEU
8	S6	148	SER
8	S6	152	ASP
9	S7	11	GLN
9	S7	12	ALA
9	S7	30	SER
10	S8	40	ALA
10	S8	116	HIS
10	S8	120	THR
10	S8	146	ARG
10	S8	152	ILE
11	S9	147	MET
11	S9	150	LEU
12	C0	30	ALA
12	C0	60	SER
13	C1	145	ALA
16	C4	18	ARG
16	C4	40	ALA
16	C4	129	LYS
16	C4	132	ARG
17	C5	127	ARG
17	C5	130	ARG
18	C6	32	ASN
18	C6	142	TYR
19	C7	42	GLN
19	C7	87	GLU
20	C8	7	GLU

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Mol	Chain	Res	Type
20	C8	90	ASN
21	C9	50	ALA
22	D0	55	PRO
22	D0	118	VAL
23	D1	7	GLN
23	D1	10	GLU
23	D1	42	GLU
23	D1	82	VAL
24	D2	92	ASN
25	D3	40	SER
25	D3	41	SER
25	D3	110	LYS
26	D4	4	ALA
26	D4	5	VAL
26	D4	6	THR
26	D4	34	ASN
26	D4	134	ALA
28	D6	63	ALA
28	D6	85	ARG
29	D7	75	GLU
33	E1	83	LYS
33	E1	93	HIS
33	E1	100	LEU
33	E1	111	GLU
34	SR	16	HIS
34	SR	185	GLN
34	SR	237	GLN
35	SM	65	THR
35	SM	100	THR
39	L2	17	THR
39	L2	115	ASN
39	L2	125	ALA
40	L3	298	PHE
40	L3	351	LEU
41	L4	146	PRO
41	L4	270	SER
41	L4	272	VAL
42	L5	132	THR
43	L6	20	LYS
44	L7	206	LYS
45	L8	253	SER
46	L9	2	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
46	L9	15	GLY
46	L9	96	HIS
46	L9	137	SER
46	L9	190	ASP
48	M1	12	LEU
48	M1	24	GLY
48	M1	151	SER
49	M3	76	THR
50	M4	9	ALA
50	M4	29	ALA
52	M6	148	LYS
53	M7	160	ALA
53	M7	164	LYS
55	M9	169	ALA
56	N0	2	ALA
57	N1	124	VAL
60	N4	70	LYS
60	N4	76	VAL
60	N4	77	LYS
60	N4	80	ARG
60	N4	97	LYS
62	N6	126	LEU
63	N7	102	GLU
64	N8	24	LYS
70	O4	14	ASN
72	O6	80	PHE
75	O9	30	ARG
75	O9	46	ARG
2	s0	127	ARG
3	s1	93	GLY
3	s1	106	THR
3	s1	207	LEU
3	s1	223	PHE
4	s2	151	PRO
4	s2	182	PRO
5	s3	45	LYS
5	s3	179	GLN
5	s3	196	ARG
6	s4	20	LEU
6	s4	96	ASN
6	s4	164	LEU
6	s4	245	LYS

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Mol	Chain	Res	Type
7	s5	56	ALA
7	s5	154	ALA
9	s7	11	GLN
10	s8	122	GLY
10	s8	136	SER
10	s8	149	SER
11	s9	156	ILE
11	s9	167	ALA
12	c0	30	ALA
14	c2	39	ASP
14	c2	54	ARG
14	c2	63	VAL
14	c2	90	LYS
15	c3	108	ASP
15	c3	139	TRP
16	c4	12	GLN
16	c4	58	TYR
16	c4	97	GLY
16	c4	114	ARG
16	c4	124	ASP
16	c4	132	ARG
17	c5	12	PHE
17	c5	14	THR
18	c6	142	TYR
20	c8	7	GLU
20	c8	94	ASP
21	c9	142	GLU
23	d1	40	ASP
25	d3	131	SER
27	d5	53	GLU
29	d7	63	LEU
29	d7	75	GLU
31	d9	11	PRO
32	e0	11	ALA
32	e0	61	SER
33	e1	99	LYS
33	e1	100	LEU
33	e1	111	GLU
33	e1	128	ALA
33	e1	136	LYS
33	e1	146	SER
34	sR	96	THR

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Mol	Chain	Res	Type
39	l2	56	ALA
39	l2	69	TYR
39	l2	130	SER
39	l2	239	ALA
40	l3	155	ALA
41	l4	4	PRO
41	l4	220	ARG
41	l4	311	HIS
41	l4	330	TYR
42	l5	258	LYS
42	l5	294	ALA
43	l6	10	TYR
44	l7	163	LEU
44	l7	191	VAL
45	l8	39	ALA
45	l8	43	LYS
45	l8	119	ALA
45	l8	133	LYS
45	l8	147	LYS
45	l8	196	ALA
47	m0	187	ALA
48	m1	153	LYS
49	m3	135	ALA
49	m3	140	SER
49	m3	162	ASN
50	m4	136	ALA
51	m5	81	TYR
51	m5	183	THR
54	m8	164	ARG
56	n0	2	ALA
59	n3	107	GLY
60	n4	64	THR
60	n4	77	LYS
61	n5	47	ALA
62	n6	6	LEU
62	n6	31	LEU
62	n6	125	LYS
63	n7	103	GLN
63	n7	125	GLY
63	n7	134	LEU
64	n8	110	GLY
68	o2	5	PRO

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Mol	Chain	Res	Type
70	o4	51	LEU
70	o4	82	ALA
77	q1	23	ARG
81	p0	33	VAL
81	p0	47	GLY
83	f	17	ALA
83	f	49	THR
83	f	63	ASP
83	f	121	GLY
83	f	152	GLU
2	S0	24	LEU
2	S0	162	CYS
2	S0	193	GLN
2	S0	194	PRO
2	S0	195	TRP
3	S1	42	ASN
3	S1	54	LEU
3	S1	131	ASP
3	S1	176	VAL
3	S1	230	ALA
4	S2	109	GLY
4	S2	146	THR
5	S3	211	PRO
6	S4	200	ARG
7	S5	45	LYS
7	S5	65	ARG
8	S6	69	LEU
9	S7	29	ASN
9	S7	75	THR
9	S7	98	ILE
9	S7	133	THR
10	S8	9	HIS
13	C1	30	ARG
14	C2	85	LYS
14	C2	115	VAL
15	C3	28	LEU
16	C4	114	ARG
17	C5	17	TYR
17	C5	53	PRO
18	C6	39	VAL
18	C6	59	LYS
18	C6	136	SER

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Mol	Chain	Res	Type
20	C8	9	GLY
20	C8	83	ALA
20	C8	144	ARG
23	D1	15	ARG
24	D2	78	ARG
25	D3	92	CYS
26	D4	53	ASP
27	D5	88	ILE
28	D6	3	LYS
28	D6	46	GLU
28	D6	47	ALA
28	D6	53	LEU
30	D8	35	ASP
33	E1	85	TYR
33	E1	97	LYS
33	E1	148	TYR
34	SR	17	ASN
34	SR	163	ASP
34	SR	194	GLY
35	SM	42	ALA
35	SM	53	ARG
35	SM	86	ASN
39	L2	55	GLY
39	L2	246	LEU
39	L2	251	LYS
40	L3	258	ALA
41	L4	5	GLN
41	L4	14	GLU
41	L4	90	PHE
41	L4	219	LEU
41	L4	233	LEU
41	L4	292	SER
42	L5	214	ASP
42	L5	259	LYS
42	L5	274	GLN
43	L6	131	LYS
46	L9	110	LYS
46	L9	117	PHE
46	L9	188	THR
46	L9	189	GLU
47	M0	71	CYS
47	M0	189	GLU

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Mol	Chain	Res	Type
48	M1	23	VAL
48	M1	64	LYS
48	M1	115	LYS
49	M3	19	GLN
49	M3	140	SER
51	M5	94	TYR
52	M6	63	ALA
54	M8	99	THR
54	M8	162	ALA
54	M8	183	GLY
55	M9	130	ASN
57	N1	18	ASP
58	N2	11	ILE
58	N2	60	GLY
59	N3	82	ALA
60	N4	68	ALA
60	N4	86	SER
63	N7	33	SER
63	N7	37	PRO
64	N8	77	LYS
67	O1	83	GLU
68	O2	70	GLY
71	O5	10	ARG
71	O5	14	LYS
71	O5	80	LEU
71	O5	119	LYS
72	O6	21	THR
72	O6	98	ARG
73	O7	18	LEU
73	O7	85	LYS
74	O8	7	ASP
79	Q3	65	ALA
3	s1	157	GLN
4	s2	88	LYS
4	s2	235	LEU
4	s2	238	SER
5	s3	44	THR
5	s3	93	ASP
6	s4	12	LEU
6	s4	51	ARG
7	s5	29	ILE
7	s5	55	ASP

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Mol	Chain	Res	Type
7	s5	71	ALA
7	s5	100	ASN
8	s6	149	LYS
8	s6	154	ARG
8	s6	216	LEU
9	s7	64	VAL
9	s7	133	THR
10	s8	78	ILE
11	s9	118	LEU
11	s9	147	MET
12	c0	3	MET
12	c0	31	LYS
13	c1	128	CYS
14	c2	87	PRO
14	c2	101	ALA
16	c4	11	SER
16	c4	37	GLU
16	c4	50	ALA
17	c5	126	VAL
17	c5	128	HIS
19	c7	68	GLY
19	c7	117	LEU
20	c8	143	ARG
21	c9	51	GLU
23	d1	2	GLU
23	d1	9	VAL
28	d6	15	ARG
28	d6	82	ARG
29	d7	29	ARG
30	d8	16	LEU
32	e0	13	LYS
33	e1	131	PHE
33	e1	137	ASP
34	sR	149	ASP
34	sR	186	PHE
34	sR	271	VAL
35	sM	43	ASP
35	sM	60	ALA
35	sM	63	ASP
39	l2	80	GLU
39	l2	247	ARG
40	l3	3	HIS

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Mol	Chain	Res	Type
40	l3	175	LYS
40	l3	348	ARG
42	l5	274	GLN
43	l6	20	LYS
44	l7	202	LEU
44	l7	229	PHE
45	l8	118	GLU
46	l9	2	LYS
46	l9	38	LEU
47	m0	195	ALA
48	m1	12	LEU
49	m3	60	ALA
49	m3	93	ILE
49	m3	152	THR
51	m5	68	ARG
59	n3	106	LYS
59	n3	109	MET
60	n4	132	GLY
61	n5	38	LEU
61	n5	45	LYS
61	n5	62	VAL
64	n8	24	LYS
65	n9	6	ASN
67	o1	90	PHE
72	o6	99	ARG
74	o8	34	ALA
78	q2	104	LEU
81	p0	102	SER
83	f	13	ALA
83	f	113	VAL
83	f	132	LYS
2	S0	109	ASN
2	S0	126	PRO
3	S1	210	ILE
3	S1	224	ASP
4	S2	39	THR
4	S2	150	GLN
4	S2	223	GLY
4	S2	235	LEU
4	S2	239	PRO
9	S7	38	LEU
9	S7	166	LEU

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Mol	Chain	Res	Type
10	S8	107	THR
11	S9	16	LYS
12	C0	25	LYS
13	C1	3	THR
15	C3	27	LYS
15	C3	68	GLY
17	C5	52	LYS
17	C5	126	VAL
19	C7	23	LYS
20	C8	36	LYS
23	D1	12	TYR
26	D4	95	GLY
27	D5	94	LYS
28	D6	64	LEU
28	D6	80	HIS
34	SR	3	SER
34	SR	139	GLN
39	L2	234	LYS
41	L4	140	HIS
41	L4	182	LEU
42	L5	81	HIS
42	L5	93	THR
42	L5	125	VAL
42	L5	181	PRO
42	L5	273	ARG
44	L7	176	TYR
44	L7	178	ILE
45	L8	36	ILE
45	L8	79	GLN
46	L9	14	GLU
46	L9	74	LEU
47	M0	18	PRO
48	M1	108	GLU
48	M1	173	ASP
49	M3	77	LEU
52	M6	16	VAL
55	M9	55	VAL
57	N1	121	ALA
57	N1	123	GLY
59	N3	109	MET
63	N7	103	GLN
65	N9	25	LYS

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Mol	Chain	Res	Type
66	O0	98	SER
67	O1	7	VAL
68	O2	68	PRO
70	O4	82	ALA
71	O5	98	SER
71	O5	99	GLN
76	Q0	79	GLU
79	Q3	7	LYS
79	Q3	45	LYS
79	Q3	51	ALA
2	s0	10	THR
2	s0	103	THR
2	s0	185	ARG
3	s1	221	PRO
4	s2	62	PRO
4	s2	150	GLN
4	s2	163	GLY
5	s3	160	SER
6	s4	30	ARG
6	s4	90	ILE
6	s4	214	LEU
7	s5	37	GLN
8	s6	65	GLN
8	s6	70	PRO
8	s6	138	ALA
11	s9	168	ARG
12	c0	24	LYS
12	c0	35	ILE
14	c2	45	LEU
14	c2	115	VAL
15	c3	29	SER
17	c5	130	ARG
18	c6	3	ALA
19	c7	113	LEU
22	d0	49	ASN
25	d3	101	GLU
26	d4	58	PHE
28	d6	8	ASN
34	sR	234	LEU
39	l2	115	ASN
39	l2	125	ALA
39	l2	241	ARG

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Mol	Chain	Res	Type
40	l3	23	ALA
41	l4	328	ASN
43	l6	147	ALA
45	l8	122	LYS
45	l8	146	LYS
45	l8	182	GLY
46	l9	189	GLU
47	m0	6	ALA
47	m0	176	LEU
48	m1	168	ASP
49	m3	46	ILE
51	m5	181	ASN
53	m7	23	ARG
54	m8	91	ALA
55	m9	130	ASN
57	n1	127	GLN
58	n2	106	ALA
61	n5	39	LYS
64	n8	78	LEU
67	o1	82	GLU
67	o1	83	GLU
74	o8	17	ARG
78	q2	66	LYS
78	q2	76	LYS
81	p0	30	VAL
83	f	15	SER
83	f	16	SER
83	f	98	GLY
83	f	127	ALA
2	S0	94	GLY
2	S0	153	SER
4	S2	182	PRO
4	S2	228	ASN
6	S4	77	ARG
11	S9	156	ILE
14	C2	66	VAL
14	C2	82	PRO
31	D9	6	VAL
33	E1	109	ASP
35	SM	88	ARG
40	L3	345	ASN
42	L5	137	ASP

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Mol	Chain	Res	Type
44	L7	26	VAL
45	L8	116	VAL
49	M3	130	GLY
50	M4	6	ILE
55	M9	124	TYR
58	N2	68	THR
59	N3	7	GLN
60	N4	74	LYS
68	O2	62	LYS
68	O2	69	SER
72	O6	34	SER
2	s0	118	PRO
3	s1	147	ALA
3	s1	177	GLN
7	s5	151	GLY
9	s7	61	PHE
9	s7	145	GLY
11	s9	185	GLY
13	c1	7	VAL
15	c3	22	ALA
16	c4	131	GLY
17	c5	52	LYS
17	c5	68	PRO
20	c8	79	TYR
21	c9	34	VAL
22	d0	96	PRO
22	d0	118	VAL
32	e0	51	ASN
34	sR	15	GLY
34	sR	105	GLY
39	l2	174	ARG
40	l3	188	ILE
41	l4	5	GLN
44	l7	56	GLU
44	l7	178	ILE
47	m0	150	GLU
64	n8	15	VAL
70	o4	14	ASN
70	o4	59	PRO
78	q2	18	ARG
83	f	97	ASP
8	S6	165	GLY

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Mol	Chain	Res	Type
9	S7	73	VAL
11	S9	15	PRO
14	C2	87	PRO
16	C4	131	GLY
25	D3	4	GLY
44	L7	217	PRO
45	L8	135	GLY
54	M8	84	VAL
58	N2	19	VAL
71	O5	77	PRO
7	s5	152	GLY
13	c1	130	PRO
24	d2	6	VAL
33	e1	112	GLY
39	l2	212	GLY
40	l3	40	PRO
47	m0	179	PRO
48	m1	120	ILE
68	o2	124	GLY
74	o8	35	GLY
3	S1	114	VAL
9	S7	109	VAL
9	S7	131	PHE
23	D1	14	PRO
40	L3	317	ILE
8	s6	165	GLY
8	s6	175	ILE
10	s8	50	GLY
11	s9	162	SER
16	c4	122	PRO
17	c5	20	VAL
35	sM	51	ARG
42	l5	91	GLY
43	l6	45	GLY
53	m7	67	ILE
59	n3	3	GLY
59	n3	41	GLY
72	o6	9	ILE
6	S4	193	GLY
9	S7	112	ARG
20	C8	82	PRO
27	D5	87	GLY

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Mol	Chain	Res	Type
28	D6	58	VAL
28	D6	59	TYR
34	SR	49	GLY
34	SR	105	GLY
35	SM	111	GLY
45	L8	30	THR
53	M7	67	ILE
56	N0	135	VAL
65	N9	21	ILE
4	s2	236	PRO
13	c1	140	VAL
26	d4	30	PRO
30	d8	30	VAL
33	e1	127	GLY
35	sM	56	GLY
52	m6	16	VAL
64	n8	28	HIS
65	n9	30	PRO
72	o6	3	VAL
7	S5	51	VAL
11	S9	18	PRO
13	C1	113	PRO
40	L3	96	PRO
41	L4	23	PRO
42	L5	170	GLY
2	s0	82	GLY
9	s7	8	ILE
13	c1	76	VAL
48	m1	117	ASP
19	C7	38	ILE
64	N8	148	ILE
13	c1	129	ARG
14	c2	102	GLY
15	c3	71	ILE
18	c6	4	VAL
21	c9	26	GLY
28	d6	20	PRO
64	n8	148	ILE
40	L3	257	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	S0	164/209 (78%)	143 (87%)	21 (13%)	5	23
2	s0	165/209 (79%)	140 (85%)	25 (15%)	3	16
3	S1	191/223 (86%)	169 (88%)	22 (12%)	7	29
3	s1	192/223 (86%)	164 (85%)	28 (15%)	4	18
4	S2	176/204 (86%)	155 (88%)	21 (12%)	6	27
4	s2	176/204 (86%)	141 (80%)	35 (20%)	1	7
5	S3	182/194 (94%)	157 (86%)	25 (14%)	4	20
5	s3	182/194 (94%)	162 (89%)	20 (11%)	8	31
6	S4	221/221 (100%)	192 (87%)	29 (13%)	5	22
6	s4	221/221 (100%)	188 (85%)	33 (15%)	4	17
7	S5	173/190 (91%)	157 (91%)	16 (9%)	11	40
7	s5	173/190 (91%)	147 (85%)	26 (15%)	3	17
8	S6	188/201 (94%)	165 (88%)	23 (12%)	6	26
8	s6	187/201 (93%)	165 (88%)	22 (12%)	6	27
9	S7	165/169 (98%)	145 (88%)	20 (12%)	6	26
9	s7	165/169 (98%)	150 (91%)	15 (9%)	12	41
10	S8	150/161 (93%)	136 (91%)	14 (9%)	11	40
10	s8	150/161 (93%)	131 (87%)	19 (13%)	5	24
11	S9	158/165 (96%)	133 (84%)	25 (16%)	3	15
11	s9	158/165 (96%)	132 (84%)	26 (16%)	3	14
12	C0	77/77 (100%)	71 (92%)	6 (8%)	16	50
12	c0	73/77 (95%)	66 (90%)	7 (10%)	10	38
13	C1	129/129 (100%)	117 (91%)	12 (9%)	11	40
13	c1	129/129 (100%)	113 (88%)	16 (12%)	6	25
14	C2	88/106 (83%)	73 (83%)	15 (17%)	2	12
14	c2	88/106 (83%)	80 (91%)	8 (9%)	12	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
15	C3	127/127 (100%)	110 (87%)	17 (13%)	5	21
15	c3	127/127 (100%)	110 (87%)	17 (13%)	5	21
16	C4	81/104 (78%)	67 (83%)	14 (17%)	2	11
16	c4	97/104 (93%)	78 (80%)	19 (20%)	1	7
17	C5	101/113 (89%)	91 (90%)	10 (10%)	10	37
17	c5	103/113 (91%)	89 (86%)	14 (14%)	5	21
18	C6	117/118 (99%)	103 (88%)	14 (12%)	6	27
18	c6	118/118 (100%)	102 (86%)	16 (14%)	5	21
19	C7	94/124 (76%)	79 (84%)	15 (16%)	3	14
19	c7	92/124 (74%)	82 (89%)	10 (11%)	8	32
20	C8	128/128 (100%)	104 (81%)	24 (19%)	2	8
20	c8	128/128 (100%)	109 (85%)	19 (15%)	4	17
21	C9	115/115 (100%)	99 (86%)	16 (14%)	4	20
21	c9	115/115 (100%)	98 (85%)	17 (15%)	4	17
22	D0	100/113 (88%)	91 (91%)	9 (9%)	12	42
22	d0	103/113 (91%)	91 (88%)	12 (12%)	7	28
23	D1	74/74 (100%)	65 (88%)	9 (12%)	6	26
23	d1	74/74 (100%)	66 (89%)	8 (11%)	8	32
24	D2	110/110 (100%)	94 (86%)	16 (14%)	4	18
24	d2	110/110 (100%)	93 (84%)	17 (16%)	3	16
25	D3	119/119 (100%)	105 (88%)	14 (12%)	6	27
25	d3	119/119 (100%)	109 (92%)	10 (8%)	14	46
26	D4	112/112 (100%)	103 (92%)	9 (8%)	15	49
26	d4	112/112 (100%)	97 (87%)	15 (13%)	5	21
27	D5	61/88 (69%)	48 (79%)	13 (21%)	1	5
27	d5	61/88 (69%)	51 (84%)	10 (16%)	3	14
28	D6	83/83 (100%)	73 (88%)	10 (12%)	6	27
28	d6	83/83 (100%)	70 (84%)	13 (16%)	3	15
29	D7	70/70 (100%)	62 (89%)	8 (11%)	7	29
29	d7	70/70 (100%)	62 (89%)	8 (11%)	7	29
30	D8	56/59 (95%)	48 (86%)	8 (14%)	4	19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
30	d8	56/59 (95%)	46 (82%)	10 (18%)	2	10
31	D9	47/48 (98%)	39 (83%)	8 (17%)	2	12
31	d9	47/48 (98%)	37 (79%)	10 (21%)	1	5
32	E0	51/53 (96%)	47 (92%)	4 (8%)	16	50
32	e0	53/53 (100%)	39 (74%)	14 (26%)	0	2
33	E1	62/66 (94%)	51 (82%)	11 (18%)	2	10
33	e1	66/66 (100%)	55 (83%)	11 (17%)	3	13
34	SR	259/261 (99%)	227 (88%)	32 (12%)	6	25
34	sR	260/261 (100%)	243 (94%)	17 (6%)	21	59
35	SM	97/183 (53%)	79 (81%)	18 (19%)	2	9
35	sM	54/183 (30%)	46 (85%)	8 (15%)	4	17
39	L2	193/195 (99%)	166 (86%)	27 (14%)	4	20
39	l2	192/195 (98%)	161 (84%)	31 (16%)	3	14
40	L3	319/322 (99%)	267 (84%)	52 (16%)	3	14
40	l3	319/322 (99%)	261 (82%)	58 (18%)	2	9
41	L4	288/288 (100%)	252 (88%)	36 (12%)	6	24
41	l4	288/288 (100%)	248 (86%)	40 (14%)	4	20
42	L5	244/244 (100%)	208 (85%)	36 (15%)	4	17
42	l5	243/244 (100%)	200 (82%)	43 (18%)	2	10
43	L6	134/152 (88%)	116 (87%)	18 (13%)	5	21
43	l6	135/152 (89%)	115 (85%)	20 (15%)	4	17
44	L7	186/204 (91%)	162 (87%)	24 (13%)	5	23
44	l7	187/204 (92%)	164 (88%)	23 (12%)	6	25
45	L8	187/207 (90%)	165 (88%)	22 (12%)	6	27
45	l8	177/207 (86%)	153 (86%)	24 (14%)	5	21
46	L9	171/171 (100%)	135 (79%)	36 (21%)	1	6
46	l9	171/171 (100%)	143 (84%)	28 (16%)	3	14
47	M0	177/186 (95%)	144 (81%)	33 (19%)	2	9
47	m0	179/186 (96%)	149 (83%)	30 (17%)	2	13
48	M1	147/150 (98%)	117 (80%)	30 (20%)	1	6
48	m1	147/150 (98%)	125 (85%)	22 (15%)	3	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
49	M3	154/158 (98%)	133 (86%)	21 (14%)	5	21
49	m3	154/158 (98%)	127 (82%)	27 (18%)	2	11
50	M4	107/108 (99%)	93 (87%)	14 (13%)	5	22
50	m4	108/108 (100%)	94 (87%)	14 (13%)	5	22
51	M5	175/175 (100%)	151 (86%)	24 (14%)	4	20
51	m5	175/175 (100%)	154 (88%)	21 (12%)	6	27
52	M6	160/161 (99%)	134 (84%)	26 (16%)	3	14
52	m6	160/161 (99%)	139 (87%)	21 (13%)	5	22
53	M7	140/145 (97%)	120 (86%)	20 (14%)	4	19
53	m7	125/145 (86%)	106 (85%)	19 (15%)	3	16
54	M8	150/150 (100%)	132 (88%)	18 (12%)	6	27
54	m8	150/150 (100%)	126 (84%)	24 (16%)	3	14
55	M9	153/153 (100%)	136 (89%)	17 (11%)	8	31
55	m9	153/153 (100%)	133 (87%)	20 (13%)	5	22
56	N0	156/156 (100%)	132 (85%)	24 (15%)	3	16
56	n0	156/156 (100%)	133 (85%)	23 (15%)	4	18
57	N1	136/136 (100%)	108 (79%)	28 (21%)	1	6
57	n1	136/136 (100%)	113 (83%)	23 (17%)	2	12
58	N2	87/106 (82%)	79 (91%)	8 (9%)	11	40
58	n2	85/106 (80%)	73 (86%)	12 (14%)	4	20
59	N3	104/104 (100%)	90 (86%)	14 (14%)	5	21
59	n3	104/104 (100%)	91 (88%)	13 (12%)	6	24
60	N4	57/129 (44%)	52 (91%)	5 (9%)	12	44
60	n4	100/129 (78%)	95 (95%)	5 (5%)	30	68
61	N5	104/117 (89%)	83 (80%)	21 (20%)	1	6
61	n5	104/117 (89%)	90 (86%)	14 (14%)	5	21
62	N6	109/109 (100%)	90 (83%)	19 (17%)	2	11
62	n6	109/109 (100%)	93 (85%)	16 (15%)	4	18
63	N7	115/115 (100%)	99 (86%)	16 (14%)	4	20
63	n7	115/115 (100%)	97 (84%)	18 (16%)	3	15
64	N8	118/118 (100%)	98 (83%)	20 (17%)	2	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
64	n8	118/118 (100%)	101 (86%)	17 (14%)	4	19
65	N9	46/46 (100%)	40 (87%)	6 (13%)	5	22
65	n9	46/46 (100%)	41 (89%)	5 (11%)	8	32
66	O0	81/87 (93%)	69 (85%)	12 (15%)	4	17
66	o0	84/87 (97%)	72 (86%)	12 (14%)	4	19
67	O1	92/96 (96%)	80 (87%)	12 (13%)	5	22
67	o1	94/96 (98%)	76 (81%)	18 (19%)	2	8
68	O2	109/110 (99%)	94 (86%)	15 (14%)	4	20
68	o2	109/110 (99%)	87 (80%)	22 (20%)	1	6
69	O3	90/90 (100%)	75 (83%)	15 (17%)	3	13
69	o3	90/90 (100%)	76 (84%)	14 (16%)	3	16
70	O4	95/102 (93%)	84 (88%)	11 (12%)	7	29
70	o4	95/102 (93%)	86 (90%)	9 (10%)	11	38
71	O5	104/104 (100%)	95 (91%)	9 (9%)	13	45
71	o5	103/104 (99%)	90 (87%)	13 (13%)	5	24
72	O6	81/81 (100%)	65 (80%)	16 (20%)	1	7
72	o6	80/81 (99%)	59 (74%)	21 (26%)	0	2
73	O7	70/70 (100%)	59 (84%)	11 (16%)	3	15
73	o7	70/70 (100%)	62 (89%)	8 (11%)	7	29
74	O8	68/68 (100%)	56 (82%)	12 (18%)	2	10
74	o8	67/68 (98%)	57 (85%)	10 (15%)	4	17
75	O9	45/45 (100%)	38 (84%)	7 (16%)	3	16
75	o9	45/45 (100%)	42 (93%)	3 (7%)	20	58
76	Q0	47/47 (100%)	37 (79%)	10 (21%)	1	5
76	q0	47/47 (100%)	39 (83%)	8 (17%)	2	12
77	Q1	23/23 (100%)	20 (87%)	3 (13%)	5	22
77	q1	23/23 (100%)	17 (74%)	6 (26%)	0	2
78	Q2	90/90 (100%)	74 (82%)	16 (18%)	2	10
78	q2	90/90 (100%)	72 (80%)	18 (20%)	1	6
79	Q3	71/71 (100%)	60 (84%)	11 (16%)	3	16
79	q3	71/71 (100%)	61 (86%)	10 (14%)	4	20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
81	p0	105/232 (45%)	93 (89%)	12 (11%)	7	29
83	f	124/133 (93%)	107 (86%)	17 (14%)	4	20
84	B	2/2 (100%)	2 (100%)	0	100	100
84	C	2/2 (100%)	2 (100%)	0	100	100
All	All	18853/20181 (93%)	16183 (86%)	2670 (14%)	4	19

All (2670) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	S0	7	PHE
2	S0	9	LEU
2	S0	23	HIS
2	S0	32	HIS
2	S0	37	VAL
2	S0	43	ASP
2	S0	52	LYS
2	S0	62	ARG
2	S0	84	ARG
2	S0	88	LYS
2	S0	96	THR
2	S0	103	THR
2	S0	111	ILE
2	S0	129	ASP
2	S0	146	LEU
2	S0	153	SER
2	S0	168	HIS
2	S0	170	ILE
2	S0	172	LEU
2	S0	188	LEU
2	S0	200	ASP
3	S1	21	VAL
3	S1	29	TRP
3	S1	30	PHE
3	S1	39	GLU
3	S1	61	LEU
3	S1	70	LEU
3	S1	81	PHE
3	S1	83	LYS
3	S1	105	PHE
3	S1	111	ARG
3	S1	180	THR

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Mol	Chain	Res	Type
3	S1	181	LEU
3	S1	184	LEU
3	S1	193	ILE
3	S1	194	ASN
3	S1	202	LYS
3	S1	203	ASP
3	S1	214	LYS
3	S1	215	VAL
3	S1	218	LEU
3	S1	222	LYS
3	S1	223	PHE
4	S2	41	LEU
4	S2	54	GLU
4	S2	69	ILE
4	S2	74	PRO
4	S2	77	GLN
4	S2	86	VAL
4	S2	89	GLN
4	S2	94	GLN
4	S2	95	ARG
4	S2	96	THR
4	S2	97	ARG
4	S2	99	LYS
4	S2	111	VAL
4	S2	137	ILE
4	S2	146	THR
4	S2	148	LEU
4	S2	159	THR
4	S2	207	LEU
4	S2	208	GLU
4	S2	222	TYR
4	S2	226	THR
5	S3	4	LEU
5	S3	5	ILE
5	S3	23	GLU
5	S3	45	LYS
5	S3	57	ASP
5	S3	65	ARG
5	S3	66	ILE
5	S3	76	ARG
5	S3	83	THR
5	S3	84	ILE

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Mol	Chain	Res	Type
5	S3	91	VAL
5	S3	92	GLN
5	S3	93	ASP
5	S3	111	ASN
5	S3	124	ARG
5	S3	158	ILE
5	S3	165	ASN
5	S3	172	THR
5	S3	176	LEU
5	S3	181	VAL
5	S3	182	LEU
5	S3	207	THR
5	S3	213	GLU
5	S3	218	LEU
5	S3	223	LYS
6	S4	3	ARG
6	S4	9	LEU
6	S4	12	LEU
6	S4	38	LEU
6	S4	45	ILE
6	S4	56	LEU
6	S4	65	LEU
6	S4	77	ARG
6	S4	92	LEU
6	S4	104	ASP
6	S4	109	PHE
6	S4	115	THR
6	S4	131	LEU
6	S4	148	ARG
6	S4	180	LEU
6	S4	181	VAL
6	S4	182	TYR
6	S4	187	ARG
6	S4	189	LEU
6	S4	192	ILE
6	S4	202	ASP
6	S4	206	ASP
6	S4	215	ASP
6	S4	220	THR
6	S4	222	LEU
6	S4	227	VAL
6	S4	240	LYS

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Mol	Chain	Res	Type
6	S4	246	LEU
6	S4	258	GLN
7	S5	25	LEU
7	S5	32	GLU
7	S5	43	PHE
7	S5	45	LYS
7	S5	46	TRP
7	S5	48	PHE
7	S5	50	GLU
7	S5	65	ARG
7	S5	79	ASN
7	S5	93	LEU
7	S5	119	ASP
7	S5	149	VAL
7	S5	156	ARG
7	S5	157	ARG
7	S5	160	VAL
7	S5	184	PHE
8	S6	7	TYR
8	S6	13	GLN
8	S6	15	THR
8	S6	25	ARG
8	S6	44	GLU
8	S6	72	ARG
8	S6	89	ASP
8	S6	109	LEU
8	S6	120	GLU
8	S6	127	THR
8	S6	128	THR
8	S6	129	VAL
8	S6	131	LYS
8	S6	151	ASP
8	S6	154	ARG
8	S6	169	TYR
8	S6	170	THR
8	S6	174	LYS
8	S6	177	ARG
8	S6	193	LEU
8	S6	211	LEU
8	S6	216	LEU
8	S6	223	LYS
9	S7	9	LEU

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Mol	Chain	Res	Type
9	S7	15	GLU
9	S7	24	PHE
9	S7	28	GLU
9	S7	38	LEU
9	S7	42	GLN
9	S7	46	ILE
9	S7	67	LEU
9	S7	77	LEU
9	S7	85	PHE
9	S7	97	ARG
9	S7	109	VAL
9	S7	110	GLN
9	S7	114	ARG
9	S7	118	LEU
9	S7	123	ASP
9	S7	126	LEU
9	S7	130	VAL
9	S7	182	VAL
9	S7	185	ILE
10	S8	5	ARG
10	S8	8	ARG
10	S8	14	THR
10	S8	21	PHE
10	S8	29	LEU
10	S8	32	GLN
10	S8	45	SER
10	S8	46	VAL
10	S8	56	ARG
10	S8	58	LEU
10	S8	62	THR
10	S8	137	LYS
10	S8	140	GLU
10	S8	187	GLU
11	S9	3	ARG
11	S9	7	THR
11	S9	21	SER
11	S9	28	LEU
11	S9	38	ASN
11	S9	39	LYS
11	S9	54	ARG
11	S9	60	LEU
11	S9	89	ASP

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Mol	Chain	Res	Type
11	S9	92	LYS
11	S9	93	LEU
11	S9	95	TYR
11	S9	97	LEU
11	S9	99	LEU
11	S9	109	LEU
11	S9	110	GLN
11	S9	112	GLN
11	S9	134	ILE
11	S9	138	LYS
11	S9	149	ARG
11	S9	151	ASP
11	S9	171	ARG
11	S9	172	VAL
11	S9	175	ARG
11	S9	182	GLU
12	C0	1	MET
12	C0	20	VAL
12	C0	76	LEU
12	C0	78	GLU
12	C0	81	ASN
12	C0	82	LEU
13	C1	9	SER
13	C1	21	ASN
13	C1	30	ARG
13	C1	40	LEU
13	C1	44	THR
13	C1	67	ARG
13	C1	69	LYS
13	C1	74	THR
13	C1	87	ARG
13	C1	123	VAL
13	C1	131	ILE
13	C1	136	ARG
14	C2	28	LEU
14	C2	33	ARG
14	C2	36	LEU
14	C2	37	VAL
14	C2	38	HIS
14	C2	43	ARG
14	C2	45	LEU
14	C2	71	ILE

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Mol	Chain	Res	Type
14	C2	74	LEU
14	C2	86	VAL
14	C2	93	ASP
14	C2	103	LEU
14	C2	126	TRP
14	C2	132	GLU
14	C2	139	HIS
15	C3	3	ARG
15	C3	9	LYS
15	C3	16	ILE
15	C3	27	LYS
15	C3	32	SER
15	C3	39	LYS
15	C3	42	ARG
15	C3	64	ARG
15	C3	66	ILE
15	C3	80	LEU
15	C3	84	ILE
15	C3	102	LEU
15	C3	105	ASN
15	C3	110	ASP
15	C3	115	LEU
15	C3	125	LEU
15	C3	140	LYS
16	C4	13	VAL
16	C4	16	VAL
16	C4	26	THR
16	C4	29	HIS
16	C4	39	ILE
16	C4	42	VAL
16	C4	81	VAL
16	C4	99	GLN
16	C4	103	ARG
16	C4	115	ILE
16	C4	125	SER
16	C4	129	LYS
16	C4	136	ARG
16	C4	137	LEU
17	C5	11	VAL
17	C5	20	VAL
17	C5	22	LEU
17	C5	32	ASP

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Mol	Chain	Res	Type
17	C5	35	LYS
17	C5	44	ARG
17	C5	52	LYS
17	C5	69	GLU
17	C5	80	MET
17	C5	121	ILE
18	C6	29	ILE
18	C6	37	THR
18	C6	39	VAL
18	C6	42	GLU
18	C6	57	LEU
18	C6	59	LYS
18	C6	66	ARG
18	C6	68	ARG
18	C6	114	ARG
18	C6	117	LEU
18	C6	123	ARG
18	C6	128	LYS
18	C6	137	ARG
18	C6	143	ARG
19	C7	3	ARG
19	C7	6	THR
19	C7	7	LYS
19	C7	25	THR
19	C7	30	THR
19	C7	38	ILE
19	C7	46	LEU
19	C7	49	LYS
19	C7	69	ILE
19	C7	84	TYR
19	C7	88	VAL
19	C7	104	ASN
19	C7	105	GLN
19	C7	112	SER
19	C7	113	LEU
20	C8	3	LEU
20	C8	5	VAL
20	C8	11	PHE
20	C8	13	HIS
20	C8	14	ILE
20	C8	17	LEU
20	C8	18	LEU

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Mol	Chain	Res	Type
20	C8	19	ASN
20	C8	28	ILE
20	C8	32	LEU
20	C8	40	ARG
20	C8	54	LEU
20	C8	57	ARG
20	C8	71	GLN
20	C8	80	LYS
20	C8	86	LEU
20	C8	92	ILE
20	C8	93	THR
20	C8	97	ASP
20	C8	108	LYS
20	C8	115	ARG
20	C8	123	ARG
20	C8	136	GLN
20	C8	138	THR
21	C9	6	VAL
21	C9	18	TYR
21	C9	22	LEU
21	C9	28	LEU
21	C9	34	VAL
21	C9	35	ASP
21	C9	54	PHE
21	C9	64	HIS
21	C9	67	MET
21	C9	68	ARG
21	C9	70	GLN
21	C9	88	VAL
21	C9	114	VAL
21	C9	125	SER
21	C9	130	ARG
21	C9	144	GLU
22	D0	18	GLN
22	D0	20	ILE
22	D0	51	VAL
22	D0	67	THR
22	D0	74	GLU
22	D0	81	THR
22	D0	84	MET
22	D0	89	ARG
22	D0	101	LYS

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Mol	Chain	Res	Type
23	D1	3	ASN
23	D1	5	LYS
23	D1	11	LEU
23	D1	32	VAL
23	D1	41	GLU
23	D1	44	ARG
23	D1	62	ARG
23	D1	78	LEU
23	D1	87	ARG
24	D2	4	SER
24	D2	24	GLN
24	D2	27	ILE
24	D2	53	ILE
24	D2	56	HIS
24	D2	65	LEU
24	D2	74	VAL
24	D2	83	ILE
24	D2	86	ILE
24	D2	93	LEU
24	D2	98	GLN
24	D2	103	ILE
24	D2	104	LEU
24	D2	120	HIS
24	D2	121	VAL
24	D2	129	VAL
25	D3	7	ARG
25	D3	9	LEU
25	D3	16	ARG
25	D3	19	ARG
25	D3	23	ARG
25	D3	40	SER
25	D3	47	SER
25	D3	65	ASN
25	D3	69	ARG
25	D3	79	ASN
25	D3	84	THR
25	D3	103	LEU
25	D3	107	PHE
25	D3	127	VAL
26	D4	34	ASN
26	D4	51	GLU
26	D4	57	VAL

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Mol	Chain	Res	Type
26	D4	62	THR
26	D4	84	LYS
26	D4	102	LYS
26	D4	123	LYS
26	D4	124	ARG
26	D4	132	ARG
27	D5	37	GLN
27	D5	40	VAL
27	D5	42	LEU
27	D5	60	VAL
27	D5	62	VAL
27	D5	67	ASP
27	D5	75	LEU
27	D5	78	ILE
27	D5	95	HIS
27	D5	96	SER
27	D5	97	LYS
27	D5	98	GLN
27	D5	102	THR
28	D6	12	LYS
28	D6	37	LYS
28	D6	38	ARG
28	D6	39	MET
28	D6	61	GLU
28	D6	64	LEU
28	D6	68	TYR
28	D6	69	ASN
28	D6	71	LEU
28	D6	91	ASP
29	D7	3	LEU
29	D7	17	ARG
29	D7	23	THR
29	D7	26	GLN
29	D7	33	LEU
29	D7	37	CYS
29	D7	42	ASN
29	D7	67	THR
30	D8	13	ILE
30	D8	14	LYS
30	D8	19	THR
30	D8	32	PHE
30	D8	39	THR

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Mol	Chain	Res	Type
30	D8	44	VAL
30	D8	58	GLU
30	D8	61	ARG
31	D9	6	VAL
31	D9	7	TRP
31	D9	12	ARG
31	D9	14	TYR
31	D9	19	ARG
31	D9	30	LEU
31	D9	36	LEU
31	D9	49	ASP
32	E0	5	HIS
32	E0	21	VAL
32	E0	31	LYS
32	E0	54	ARG
33	E1	85	TYR
33	E1	89	LYS
33	E1	98	VAL
33	E1	100	LEU
33	E1	102	VAL
33	E1	103	LEU
33	E1	108	VAL
33	E1	130	VAL
33	E1	135	HIS
33	E1	147	VAL
33	E1	151	ASN
34	SR	6	VAL
34	SR	17	ASN
34	SR	31	ASN
34	SR	39	ASP
34	SR	45	TRP
34	SR	52	GLN
34	SR	74	THR
34	SR	76	ASP
34	SR	81	LEU
34	SR	106	HIS
34	SR	116	ASP
34	SR	117	LYS
34	SR	131	ILE
34	SR	136	ILE
34	SR	137	LYS
34	SR	140	CYS

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Mol	Chain	Res	Type
34	SR	159	ASN
34	SR	184	ASN
34	SR	223	TRP
34	SR	229	LYS
34	SR	234	LEU
34	SR	238	ASP
34	SR	242	SER
34	SR	264	SER
34	SR	268	GLN
34	SR	269	TYR
34	SR	275	ARG
34	SR	290	VAL
34	SR	292	LEU
34	SR	308	ASN
34	SR	316	MET
34	SR	317	THR
35	SM	33	LYS
35	SM	34	LYS
35	SM	46	LYS
35	SM	48	ARG
35	SM	51	ARG
35	SM	53	ARG
35	SM	68	ARG
35	SM	70	ASN
35	SM	72	ARG
35	SM	77	THR
35	SM	84	LYS
35	SM	91	THR
35	SM	94	HIS
35	SM	100	THR
35	SM	116	GLU
35	SM	117	LEU
35	SM	131	ILE
35	SM	140	ASP
39	L2	14	SER
39	L2	20	THR
39	L2	28	LYS
39	L2	32	LEU
39	L2	44	ILE
39	L2	45	VAL
39	L2	70	ARG
39	L2	72	ARG

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Mol	Chain	Res	Type
39	L2	82	VAL
39	L2	84	THR
39	L2	101	VAL
39	L2	111	THR
39	L2	113	VAL
39	L2	116	VAL
39	L2	130	SER
39	L2	157	VAL
39	L2	165	VAL
39	L2	179	LEU
39	L2	180	LEU
39	L2	192	LYS
39	L2	199	THR
39	L2	204	MET
39	L2	207	VAL
39	L2	208	ASP
39	L2	227	ARG
39	L2	238	ILE
39	L2	252	THR
40	L3	7	GLU
40	L3	17	LEU
40	L3	19	ARG
40	L3	20	LYS
40	L3	21	ARG
40	L3	25	ILE
40	L3	30	LYS
40	L3	38	SER
40	L3	55	THR
40	L3	67	PHE
40	L3	72	VAL
40	L3	76	VAL
40	L3	79	VAL
40	L3	84	VAL
40	L3	85	VAL
40	L3	97	ARG
40	L3	103	THR
40	L3	104	THR
40	L3	114	VAL
40	L3	116	ARG
40	L3	128	LYS
40	L3	140	ASP
40	L3	164	THR

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Mol	Chain	Res	Type
40	L3	167	ARG
40	L3	173	GLN
40	L3	183	LEU
40	L3	187	SER
40	L3	188	ILE
40	L3	196	ARG
40	L3	206	ASP
40	L3	213	GLU
40	L3	226	PHE
40	L3	235	THR
40	L3	238	LEU
40	L3	244	ARG
40	L3	246	LEU
40	L3	252	ILE
40	L3	275	ARG
40	L3	284	ARG
40	L3	293	ASN
40	L3	296	THR
40	L3	301	THR
40	L3	304	THR
40	L3	305	ILE
40	L3	320	ASP
40	L3	328	ILE
40	L3	332	ARG
40	L3	343	TYR
40	L3	353	GLU
40	L3	354	VAL
40	L3	386	ASP
40	L3	387	LEU
41	L4	55	LYS
41	L4	74	ILE
41	L4	93	MET
41	L4	99	MET
41	L4	105	THR
41	L4	120	TYR
41	L4	124	SER
41	L4	133	SER
41	L4	136	LEU
41	L4	148	ILE
41	L4	153	SER
41	L4	156	LEU
41	L4	172	VAL

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Mol	Chain	Res	Type
41	L4	179	LEU
41	L4	187	LEU
41	L4	193	LYS
41	L4	194	TYR
41	L4	203	ARG
41	L4	206	LEU
41	L4	227	THR
41	L4	230	VAL
41	L4	246	ARG
41	L4	252	GLU
41	L4	258	LEU
41	L4	259	ASP
41	L4	285	ASP
41	L4	287	THR
41	L4	290	ILE
41	L4	307	GLN
41	L4	311	HIS
41	L4	313	LEU
41	L4	316	ASN
41	L4	333	VAL
41	L4	338	LYS
41	L4	346	LYS
41	L4	349	THR
42	L5	8	LYS
42	L5	14	SER
42	L5	23	ARG
42	L5	35	ARG
42	L5	41	LYS
42	L5	66	SER
42	L5	69	ILE
42	L5	70	THR
42	L5	81	HIS
42	L5	93	THR
42	L5	95	TRP
42	L5	105	ILE
42	L5	110	LEU
42	L5	115	LEU
42	L5	128	GLU
42	L5	130	GLU
42	L5	131	LEU
42	L5	137	ASP
42	L5	140	ARG

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Mol	Chain	Res	Type
42	L5	142	PHE
42	L5	146	LEU
42	L5	155	THR
42	L5	163	LEU
42	L5	177	GLU
42	L5	185	PHE
42	L5	188	GLU
42	L5	190	ILE
42	L5	194	LEU
42	L5	208	MET
42	L5	211	LEU
42	L5	222	LEU
42	L5	231	ILE
42	L5	263	GLU
42	L5	277	LEU
42	L5	279	LYS
42	L5	296	GLN
43	L6	5	LYS
43	L6	21	THR
43	L6	52	VAL
43	L6	57	HIS
43	L6	64	LEU
43	L6	65	ILE
43	L6	78	ARG
43	L6	79	VAL
43	L6	84	VAL
43	L6	89	THR
43	L6	90	LYS
43	L6	98	VAL
43	L6	134	ARG
43	L6	137	ASP
43	L6	154	LEU
43	L6	155	LEU
43	L6	160	SER
43	L6	162	SER
44	L7	24	GLU
44	L7	25	GLN
44	L7	26	VAL
44	L7	38	LYS
44	L7	60	ARG
44	L7	82	LYS
44	L7	88	ARG

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Mol	Chain	Res	Type
44	L7	93	ASN
44	L7	100	ARG
44	L7	105	LEU
44	L7	107	ARG
44	L7	110	ARG
44	L7	115	THR
44	L7	120	THR
44	L7	124	LEU
44	L7	126	LEU
44	L7	128	LYS
44	L7	151	ARG
44	L7	158	LYS
44	L7	175	LYS
44	L7	179	LEU
44	L7	182	ASP
44	L7	184	LEU
44	L7	239	LEU
45	L8	27	THR
45	L8	41	GLN
45	L8	47	SER
45	L8	65	LEU
45	L8	71	VAL
45	L8	74	THR
45	L8	79	GLN
45	L8	83	ASP
45	L8	109	LEU
45	L8	134	TYR
45	L8	136	LEU
45	L8	145	ASN
45	L8	150	LEU
45	L8	156	ASP
45	L8	160	ILE
45	L8	169	LEU
45	L8	181	LYS
45	L8	185	ARG
45	L8	190	VAL
45	L8	221	ASN
45	L8	230	LYS
45	L8	251	LYS
46	L9	1	MET
46	L9	5	GLN
46	L9	6	THR

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Mol	Chain	Res	Type
46	L9	9	GLN
46	L9	19	SER
46	L9	41	ILE
46	L9	42	ASP
46	L9	44	THR
46	L9	48	VAL
46	L9	49	ASN
46	L9	51	GLN
46	L9	52	LEU
46	L9	65	VAL
46	L9	68	LEU
46	L9	69	ARG
46	L9	70	THR
46	L9	71	VAL
46	L9	73	SER
46	L9	77	ASN
46	L9	92	TYR
46	L9	103	ILE
46	L9	118	LEU
46	L9	120	ASP
46	L9	132	VAL
46	L9	139	ASN
46	L9	149	ASN
46	L9	151	VAL
46	L9	157	ASN
46	L9	161	LEU
46	L9	162	GLN
46	L9	164	ILE
46	L9	172	ILE
46	L9	173	ARG
46	L9	177	ASP
46	L9	182	SER
46	L9	189	GLU
47	M0	3	ARG
47	M0	12	GLN
47	M0	26	VAL
47	M0	30	LYS
47	M0	31	ILE
47	M0	32	ARG
47	M0	48	LEU
47	M0	51	HIS
47	M0	52	LEU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
47	M0	58	GLU
47	M0	60	LEU
47	M0	61	SER
47	M0	63	GLU
47	M0	74	LYS
47	M0	82	ARG
47	M0	83	ASP
47	M0	87	LEU
47	M0	128	ARG
47	M0	130	ASP
47	M0	137	SER
47	M0	138	VAL
47	M0	139	ARG
47	M0	163	GLN
47	M0	165	ILE
47	M0	167	LEU
47	M0	174	THR
47	M0	177	ASP
47	M0	178	ARG
47	M0	185	ARG
47	M0	189	GLU
47	M0	191	LYS
47	M0	203	LYS
47	M0	208	ASN
48	M1	10	ARG
48	M1	12	LEU
48	M1	13	LYS
48	M1	19	LEU
48	M1	26	SER
48	M1	28	ASP
48	M1	39	GLN
48	M1	44	THR
48	M1	46	VAL
48	M1	61	ARG
48	M1	67	VAL
48	M1	80	LEU
48	M1	81	GLU
48	M1	82	ARG
48	M1	85	LYS
48	M1	95	ASN
48	M1	101	ASN
48	M1	106	ILE

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Mol	Chain	Res	Type
48	M1	107	ASP
48	M1	111	ASP
48	M1	112	LEU
48	M1	115	LYS
48	M1	137	ARG
48	M1	138	VAL
48	M1	140	ARG
48	M1	145	LYS
48	M1	155	THR
48	M1	157	GLU
48	M1	165	GLN
48	M1	166	LYS
49	M3	17	HIS
49	M3	22	VAL
49	M3	23	LYS
49	M3	24	VAL
49	M3	36	ARG
49	M3	53	LEU
49	M3	54	LEU
49	M3	55	ARG
49	M3	57	VAL
49	M3	59	ARG
49	M3	67	ARG
49	M3	104	ARG
49	M3	114	GLN
49	M3	117	LYS
49	M3	121	SER
49	M3	122	LYS
49	M3	124	ILE
49	M3	131	LYS
49	M3	134	GLU
49	M3	138	VAL
49	M3	169	THR
50	M4	4	ASP
50	M4	8	LYS
50	M4	14	LEU
50	M4	15	VAL
50	M4	20	VAL
50	M4	43	LYS
50	M4	50	LYS
50	M4	53	VAL
50	M4	64	VAL

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
50	M4	65	LEU
50	M4	69	THR
50	M4	72	LEU
50	M4	90	VAL
50	M4	91	CYS
51	M5	7	LEU
51	M5	17	ASP
51	M5	18	VAL
51	M5	22	LEU
51	M5	46	ASP
51	M5	80	THR
51	M5	85	THR
51	M5	94	TYR
51	M5	99	ARG
51	M5	106	VAL
51	M5	117	ASN
51	M5	133	ILE
51	M5	134	LEU
51	M5	138	GLN
51	M5	151	ILE
51	M5	153	ASP
51	M5	157	LYS
51	M5	167	THR
51	M5	183	THR
51	M5	184	LYS
51	M5	188	ARG
51	M5	190	THR
51	M5	196	THR
51	M5	198	SER
52	M6	4	GLU
52	M6	22	VAL
52	M6	31	GLN
52	M6	51	LYS
52	M6	58	LEU
52	M6	68	ARG
52	M6	77	SER
52	M6	78	ARG
52	M6	79	ILE
52	M6	84	LEU
52	M6	85	ARG
52	M6	94	ARG
52	M6	106	GLU

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Mol	Chain	Res	Type
52	M6	113	ASP
52	M6	116	LYS
52	M6	122	GLN
52	M6	124	LEU
52	M6	128	ARG
52	M6	143	THR
52	M6	144	SER
52	M6	149	TYR
52	M6	155	LYS
52	M6	160	ARG
52	M6	163	SER
52	M6	184	THR
52	M6	189	ASP
53	M7	7	THR
53	M7	24	VAL
53	M7	29	THR
53	M7	36	ILE
53	M7	42	THR
53	M7	52	LEU
53	M7	75	GLU
53	M7	78	VAL
53	M7	110	THR
53	M7	112	LEU
53	M7	118	GLN
53	M7	127	ARG
53	M7	129	THR
53	M7	137	ASN
53	M7	138	LYS
53	M7	142	SER
53	M7	144	SER
53	M7	148	LEU
53	M7	168	LEU
53	M7	171	ARG
54	M8	17	THR
54	M8	21	SER
54	M8	22	ASP
54	M8	24	VAL
54	M8	26	LEU
54	M8	32	LEU
54	M8	49	LEU
54	M8	69	ARG
54	M8	86	THR

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Mol	Chain	Res	Type
54	M8	105	ARG
54	M8	115	VAL
54	M8	127	LEU
54	M8	135	GLN
54	M8	136	ASN
54	M8	138	LEU
54	M8	171	LYS
54	M8	178	ARG
54	M8	180	ARG
55	M9	22	VAL
55	M9	44	LEU
55	M9	49	THR
55	M9	51	VAL
55	M9	60	LYS
55	M9	70	LYS
55	M9	74	ARG
55	M9	76	SER
55	M9	103	ARG
55	M9	106	LEU
55	M9	110	ARG
55	M9	134	HIS
55	M9	138	LEU
55	M9	139	VAL
55	M9	143	ILE
55	M9	165	LYS
55	M9	180	LYS
56	N0	1	MET
56	N0	16	THR
56	N0	34	GLU
56	N0	45	LEU
56	N0	50	LYS
56	N0	61	ILE
56	N0	71	LYS
56	N0	79	VAL
56	N0	81	TYR
56	N0	87	THR
56	N0	88	HIS
56	N0	92	LYS
56	N0	97	VAL
56	N0	100	VAL
56	N0	115	ARG
56	N0	117	ARG

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Mol	Chain	Res	Type
56	N0	122	HIS
56	N0	132	THR
56	N0	136	LYS
56	N0	138	GLN
56	N0	149	LYS
56	N0	166	LYS
56	N0	171	PHE
56	N0	172	TYR
57	N1	12	ARG
57	N1	16	GLN
57	N1	18	ASP
57	N1	25	VAL
57	N1	26	HIS
57	N1	27	LEU
57	N1	32	LYS
57	N1	55	LYS
57	N1	72	VAL
57	N1	75	ILE
57	N1	78	LYS
57	N1	79	MET
57	N1	80	VAL
57	N1	83	ARG
57	N1	88	ARG
57	N1	89	LEU
57	N1	93	VAL
57	N1	101	CYS
57	N1	103	GLN
57	N1	104	GLU
57	N1	106	LEU
57	N1	118	GLU
57	N1	124	VAL
57	N1	126	VAL
57	N1	128	LEU
57	N1	131	GLN
57	N1	139	ARG
57	N1	141	VAL
58	N2	10	LYS
58	N2	16	THR
58	N2	19	VAL
58	N2	50	LEU
58	N2	52	ASN
58	N2	66	VAL

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
58	N2	70	LYS
58	N2	100	THR
59	N3	9	THR
59	N3	44	SER
59	N3	48	ARG
59	N3	54	LEU
59	N3	69	LEU
59	N3	73	VAL
59	N3	86	ARG
59	N3	88	ARG
59	N3	98	ASN
59	N3	102	ILE
59	N3	108	GLU
59	N3	115	THR
59	N3	124	ASP
59	N3	128	ARG
60	N4	1	MET
60	N4	4	GLU
60	N4	5	ILE
60	N4	19	THR
60	N4	64	THR
61	N5	25	LYS
61	N5	27	ARG
61	N5	29	SER
61	N5	36	LYS
61	N5	37	THR
61	N5	38	LEU
61	N5	39	LYS
61	N5	63	ILE
61	N5	69	SER
61	N5	70	GLU
61	N5	77	GLU
61	N5	78	ASP
61	N5	86	VAL
61	N5	108	LEU
61	N5	113	LEU
61	N5	115	ARG
61	N5	119	THR
61	N5	125	ARG
61	N5	133	LEU
61	N5	135	ILE
61	N5	137	ASN

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Mol	Chain	Res	Type
62	N6	3	LYS
62	N6	5	SER
62	N6	9	SER
62	N6	13	ARG
62	N6	28	ARG
62	N6	37	LYS
62	N6	38	GLU
62	N6	39	LEU
62	N6	40	ARG
62	N6	42	GLN
62	N6	45	ILE
62	N6	55	GLU
62	N6	56	VAL
62	N6	57	LEU
62	N6	74	TYR
62	N6	76	LEU
62	N6	115	ARG
62	N6	122	LYS
62	N6	127	GLU
63	N7	17	ARG
63	N7	24	VAL
63	N7	30	ASP
63	N7	34	LYS
63	N7	46	ILE
63	N7	54	THR
63	N7	55	LYS
63	N7	80	LEU
63	N7	92	PHE
63	N7	94	SER
63	N7	100	THR
63	N7	102	GLU
63	N7	103	GLN
63	N7	121	ARG
63	N7	122	HIS
63	N7	135	ARG
64	N8	4	ARG
64	N8	6	THR
64	N8	8	THR
64	N8	10	LYS
64	N8	12	ARG
64	N8	34	MET
64	N8	42	ARG

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Mol	Chain	Res	Type
64	N8	46	ASP
64	N8	56	VAL
64	N8	60	TYR
64	N8	73	LEU
64	N8	76	ASP
64	N8	88	ASP
64	N8	91	LEU
64	N8	115	LYS
64	N8	118	ILE
64	N8	124	ILE
64	N8	130	VAL
64	N8	133	LEU
64	N8	135	GLU
65	N9	3	LYS
65	N9	8	THR
65	N9	13	THR
65	N9	18	ARG
65	N9	22	LYS
65	N9	59	LYS
66	O0	30	THR
66	O0	33	SER
66	O0	41	LEU
66	O0	52	ARG
66	O0	54	SER
66	O0	59	TYR
66	O0	61	MET
66	O0	67	VAL
66	O0	76	GLU
66	O0	83	LYS
66	O0	87	VAL
66	O0	100	ILE
67	O1	6	ASP
67	O1	8	VAL
67	O1	26	LYS
67	O1	31	ARG
67	O1	41	LYS
67	O1	55	LEU
67	O1	64	VAL
67	O1	73	LEU
67	O1	79	ARG
67	O1	82	GLU
67	O1	84	ASP

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
67	O1	106	THR
68	O2	3	SER
68	O2	9	ILE
68	O2	19	ARG
68	O2	31	ASN
68	O2	33	ARG
68	O2	35	GLN
68	O2	41	VAL
68	O2	54	LYS
68	O2	61	LYS
68	O2	73	THR
68	O2	95	GLU
68	O2	104	ASN
68	O2	106	VAL
68	O2	109	LEU
68	O2	126	LEU
69	O3	3	GLU
69	O3	7	LEU
69	O3	14	LEU
69	O3	15	SER
69	O3	20	LYS
69	O3	22	VAL
69	O3	24	ASN
69	O3	37	THR
69	O3	48	ARG
69	O3	59	VAL
69	O3	74	THR
69	O3	78	SER
69	O3	81	VAL
69	O3	98	VAL
69	O3	105	SER
70	O4	6	THR
70	O4	8	ARG
70	O4	23	VAL
70	O4	24	LYS
70	O4	29	ILE
70	O4	38	LEU
70	O4	58	ARG
70	O4	65	VAL
70	O4	71	THR
70	O4	102	LYS
70	O4	104	VAL

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Mol	Chain	Res	Type
71	O5	13	SER
71	O5	20	GLN
71	O5	47	VAL
71	O5	48	ARG
71	O5	49	LYS
71	O5	89	ARG
71	O5	101	THR
71	O5	102	GLU
71	O5	105	ARG
72	O6	11	LEU
72	O6	13	LYS
72	O6	21	THR
72	O6	25	LYS
72	O6	26	ILE
72	O6	28	TYR
72	O6	36	ARG
72	O6	42	SER
72	O6	57	LEU
72	O6	60	LEU
72	O6	68	ARG
72	O6	76	ARG
72	O6	80	PHE
72	O6	90	MET
72	O6	98	ARG
72	O6	99	ARG
73	O7	14	LYS
73	O7	24	ARG
73	O7	25	ARG
73	O7	31	LYS
73	O7	33	THR
73	O7	36	SER
73	O7	44	THR
73	O7	59	THR
73	O7	67	LEU
73	O7	75	LYS
73	O7	85	LYS
74	O8	8	ILE
74	O8	27	ILE
74	O8	28	ASN
74	O8	29	LYS
74	O8	32	ASN
74	O8	39	ARG

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Mol	Chain	Res	Type
74	O8	46	ARG
74	O8	53	THR
74	O8	65	LEU
74	O8	67	GLN
74	O8	69	LEU
74	O8	77	ARG
75	O9	4	GLN
75	O9	5	LYS
75	O9	21	ARG
75	O9	23	LEU
75	O9	27	ILE
75	O9	28	ARG
75	O9	29	LEU
76	Q0	79	GLU
76	Q0	81	SER
76	Q0	85	LEU
76	Q0	97	ARG
76	Q0	98	LYS
76	Q0	99	CYS
76	Q0	112	LYS
76	Q0	113	ARG
76	Q0	114	LYS
76	Q0	127	LEU
77	Q1	11	ARG
77	Q1	13	LEU
77	Q1	23	ARG
78	Q2	7	THR
78	Q2	16	THR
78	Q2	17	CYS
78	Q2	20	HIS
78	Q2	21	THR
78	Q2	26	THR
78	Q2	47	GLN
78	Q2	65	THR
78	Q2	78	LYS
78	Q2	83	LEU
78	Q2	84	THR
78	Q2	85	LEU
78	Q2	87	ARG
78	Q2	88	CYS
78	Q2	93	LEU
78	Q2	105	GLN

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Mol	Chain	Res	Type
79	Q3	5	THR
79	Q3	7	LYS
79	Q3	11	THR
79	Q3	28	LYS
79	Q3	32	GLN
79	Q3	45	LYS
79	Q3	60	CYS
79	Q3	70	THR
79	Q3	73	THR
79	Q3	82	THR
79	Q3	84	ARG
2	s0	22	THR
2	s0	31	VAL
2	s0	45	VAL
2	s0	46	HIS
2	s0	59	LEU
2	s0	72	ASP
2	s0	87	LEU
2	s0	96	THR
2	s0	108	THR
2	s0	111	ILE
2	s0	124	THR
2	s0	131	GLN
2	s0	135	GLU
2	s0	144	ILE
2	s0	151	SER
2	s0	153	SER
2	s0	154	GLU
2	s0	158	VAL
2	s0	172	LEU
2	s0	184	LEU
2	s0	188	LEU
2	s0	189	VAL
2	s0	191	ARG
2	s0	198	MET
2	s0	200	ASP
3	s1	21	VAL
3	s1	25	THR
3	s1	26	ARG
3	s1	47	LEU
3	s1	65	VAL
3	s1	66	VAL

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Mol	Chain	Res	Type
3	s1	68	VAL
3	s1	70	LEU
3	s1	74	GLN
3	s1	77	GLU
3	s1	78	ASP
3	s1	81	PHE
3	s1	101	HIS
3	s1	104	ASP
3	s1	108	ASP
3	s1	110	LEU
3	s1	124	ASN
3	s1	135	LEU
3	s1	144	ARG
3	s1	152	ARG
3	s1	153	HIS
3	s1	157	GLN
3	s1	173	THR
3	s1	181	LEU
3	s1	208	GLN
3	s1	209	ASN
3	s1	212	VAL
3	s1	216	LYS
4	s2	46	LYS
4	s2	51	THR
4	s2	53	ILE
4	s2	54	GLU
4	s2	55	GLU
4	s2	58	LEU
4	s2	69	ILE
4	s2	70	ASP
4	s2	72	LEU
4	s2	78	ASP
4	s2	82	ASN
4	s2	83	ILE
4	s2	84	LYS
4	s2	91	ARG
4	s2	94	GLN
4	s2	95	ARG
4	s2	96	THR
4	s2	97	ARG
4	s2	111	VAL
4	s2	115	ILE

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Mol	Chain	Res	Type
4	s2	125	ILE
4	s2	141	ARG
4	s2	152	HIS
4	s2	159	THR
4	s2	166	THR
4	s2	194	GLU
4	s2	207	LEU
4	s2	218	ILE
4	s2	221	THR
4	s2	222	TYR
4	s2	224	PHE
4	s2	226	THR
4	s2	228	ASN
4	s2	232	GLU
4	s2	237	VAL
5	s3	21	LEU
5	s3	26	THR
5	s3	37	VAL
5	s3	66	ILE
5	s3	67	ASN
5	s3	79	TYR
5	s3	84	ILE
5	s3	90	ARG
5	s3	91	VAL
5	s3	96	LEU
5	s3	104	SER
5	s3	116	ARG
5	s3	127	MET
5	s3	128	GLU
5	s3	158	ILE
5	s3	162	GLN
5	s3	168	ILE
5	s3	176	LEU
5	s3	212	LYS
5	s3	218	LEU
6	s4	9	LEU
6	s4	12	LEU
6	s4	23	LEU
6	s4	24	SER
6	s4	38	LEU
6	s4	42	LEU
6	s4	48	LEU

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Mol	Chain	Res	Type
6	s4	49	ARG
6	s4	51	ARG
6	s4	56	LEU
6	s4	67	GLN
6	s4	70	VAL
6	s4	75	LYS
6	s4	95	THR
6	s4	98	ASN
6	s4	104	ASP
6	s4	120	SER
6	s4	127	LYS
6	s4	128	LYS
6	s4	131	LEU
6	s4	147	ILE
6	s4	148	ARG
6	s4	163	ASP
6	s4	164	LEU
6	s4	176	ASP
6	s4	182	TYR
6	s4	191	ARG
6	s4	214	LEU
6	s4	220	THR
6	s4	222	LEU
6	s4	237	SER
6	s4	238	LEU
6	s4	252	ARG
7	s5	25	LEU
7	s5	34	GLN
7	s5	40	ILE
7	s5	60	ASP
7	s5	63	GLN
7	s5	65	ARG
7	s5	66	GLN
7	s5	68	ILE
7	s5	82	PHE
7	s5	89	ILE
7	s5	93	LEU
7	s5	98	MET
7	s5	99	MET
7	s5	102	ARG
7	s5	119	ASP
7	s5	124	LEU

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Mol	Chain	Res	Type
7	s5	146	THR
7	s5	157	ARG
7	s5	162	VAL
7	s5	189	THR
7	s5	193	THR
7	s5	194	LEU
7	s5	203	LYS
7	s5	211	ILE
7	s5	216	GLU
7	s5	222	LYS
8	s6	15	THR
8	s6	19	ASP
8	s6	21	GLU
8	s6	31	ARG
8	s6	34	GLN
8	s6	57	ASP
8	s6	64	LYS
8	s6	71	THR
8	s6	76	LEU
8	s6	78	THR
8	s6	93	LYS
8	s6	109	LEU
8	s6	119	GLN
8	s6	120	GLU
8	s6	128	THR
8	s6	129	VAL
8	s6	133	LEU
8	s6	163	THR
8	s6	179	VAL
8	s6	191	ARG
8	s6	194	LYS
8	s6	215	ARG
9	s7	9	LEU
9	s7	11	GLN
9	s7	50	ASP
9	s7	62	VAL
9	s7	75	THR
9	s7	77	LEU
9	s7	87	ASP
9	s7	97	ARG
9	s7	108	GLN
9	s7	114	ARG

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Mol	Chain	Res	Type
9	s7	143	LEU
9	s7	144	VAL
9	s7	160	GLN
9	s7	163	ASP
9	s7	176	LEU
10	s8	4	SER
10	s8	10	LYS
10	s8	18	ARG
10	s8	25	ARG
10	s8	29	LEU
10	s8	46	VAL
10	s8	58	LEU
10	s8	59	ARG
10	s8	61	GLU
10	s8	74	LYS
10	s8	76	THR
10	s8	89	GLU
10	s8	95	THR
10	s8	120	THR
10	s8	151	LYS
10	s8	152	ILE
10	s8	155	SER
10	s8	176	SER
10	s8	183	ILE
11	s9	3	ARG
11	s9	6	ARG
11	s9	7	THR
11	s9	22	SER
11	s9	28	LEU
11	s9	46	SER
11	s9	49	LEU
11	s9	54	ARG
11	s9	61	THR
11	s9	78	ARG
11	s9	82	ARG
11	s9	89	ASP
11	s9	93	LEU
11	s9	97	LEU
11	s9	103	ASP
11	s9	109	LEU
11	s9	111	THR
11	s9	120	LYS

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Mol	Chain	Res	Type
11	s9	126	ARG
11	s9	127	VAL
11	s9	132	ARG
11	s9	134	ILE
11	s9	157	ASP
11	s9	168	ARG
11	s9	172	VAL
11	s9	180	LYS
12	c0	7	ASP
12	c0	15	LEU
12	c0	27	PHE
12	c0	47	GLN
12	c0	57	THR
12	c0	67	THR
12	c0	71	GLU
13	c1	5	LEU
13	c1	10	GLU
13	c1	26	LYS
13	c1	31	THR
13	c1	33	ARG
13	c1	35	TYR
13	c1	40	LEU
13	c1	44	THR
13	c1	47	THR
13	c1	60	PHE
13	c1	67	ARG
13	c1	83	THR
13	c1	90	TYR
13	c1	111	VAL
13	c1	117	VAL
13	c1	129	ARG
14	c2	38	HIS
14	c2	45	LEU
14	c2	54	ARG
14	c2	61	VAL
14	c2	62	LEU
14	c2	71	ILE
14	c2	103	LEU
14	c2	136	ILE
15	c3	6	SER
15	c3	21	ASN
15	c3	30	SER

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Mol	Chain	Res	Type
15	c3	46	THR
15	c3	60	VAL
15	c3	66	ILE
15	c3	67	THR
15	c3	70	LYS
15	c3	74	ILE
15	c3	75	LEU
15	c3	78	ASN
15	c3	84	ILE
15	c3	101	HIS
15	c3	115	LEU
15	c3	125	LEU
15	c3	128	TYR
15	c3	149	LEU
16	c4	20	TYR
16	c4	23	PHE
16	c4	26	THR
16	c4	51	ASP
16	c4	52	ARG
16	c4	61	MET
16	c4	81	VAL
16	c4	92	LYS
16	c4	93	THR
16	c4	102	LEU
16	c4	107	ARG
16	c4	108	SER
16	c4	114	ARG
16	c4	121	VAL
16	c4	123	SER
16	c4	124	ASP
16	c4	133	ARG
16	c4	136	ARG
16	c4	137	LEU
17	c5	12	PHE
17	c5	16	SER
17	c5	24	LYS
17	c5	27	GLU
17	c5	28	MET
17	c5	36	LEU
17	c5	61	ARG
17	c5	69	GLU
17	c5	84	ILE

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Mol	Chain	Res	Type
17	c5	86	VAL
17	c5	124	THR
17	c5	126	VAL
17	c5	127	ARG
17	c5	128	HIS
18	c6	23	LYS
18	c6	26	LYS
18	c6	28	LEU
18	c6	37	THR
18	c6	43	ILE
18	c6	53	LEU
18	c6	55	VAL
18	c6	57	LEU
18	c6	68	ARG
18	c6	70	THR
18	c6	105	LEU
18	c6	114	ARG
18	c6	115	THR
18	c6	118	ILE
18	c6	127	LYS
18	c6	137	ARG
19	c7	5	ARG
19	c7	6	THR
19	c7	8	THR
19	c7	29	GLN
19	c7	34	LEU
19	c7	46	LEU
19	c7	60	ARG
19	c7	72	LYS
19	c7	112	SER
19	c7	113	LEU
20	c8	4	VAL
20	c8	5	VAL
20	c8	14	ILE
20	c8	15	LEU
20	c8	18	LEU
20	c8	28	ILE
20	c8	33	THR
20	c8	38	VAL
20	c8	40	ARG
20	c8	41	ARG
20	c8	51	ASP

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Mol	Chain	Res	Type
20	c8	85	PHE
20	c8	94	ASP
20	c8	105	VAL
20	c8	116	LEU
20	c8	120	ARG
20	c8	133	VAL
20	c8	134	ARG
20	c8	138	THR
21	c9	6	VAL
21	c9	7	ARG
21	c9	8	ASP
21	c9	9	VAL
21	c9	28	LEU
21	c9	30	VAL
21	c9	57	ARG
21	c9	68	ARG
21	c9	71	VAL
21	c9	86	ARG
21	c9	111	ILE
21	c9	123	ARG
21	c9	126	GLU
21	c9	131	ASP
21	c9	134	ARG
21	c9	140	LEU
21	c9	142	GLU
22	d0	20	ILE
22	d0	21	LYS
22	d0	25	THR
22	d0	44	ASN
22	d0	48	HIS
22	d0	57	ARG
22	d0	60	THR
22	d0	62	VAL
22	d0	70	THR
22	d0	77	LYS
22	d0	103	ILE
22	d0	115	GLU
23	d1	5	LYS
23	d1	8	LEU
23	d1	27	ASP
23	d1	32	VAL
23	d1	56	SER

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Mol	Chain	Res	Type
23	d1	62	ARG
23	d1	68	SER
23	d1	81	ASN
24	d2	7	LEU
24	d2	9	ASP
24	d2	22	LYS
24	d2	25	VAL
24	d2	37	PHE
24	d2	42	GLN
24	d2	57	ARG
24	d2	58	SER
24	d2	65	LEU
24	d2	85	ASP
24	d2	93	LEU
24	d2	98	GLN
24	d2	99	PHE
24	d2	103	ILE
24	d2	105	THR
24	d2	126	LEU
24	d2	129	VAL
25	d3	19	ARG
25	d3	28	ASN
25	d3	40	SER
25	d3	73	ARG
25	d3	75	GLN
25	d3	84	THR
25	d3	95	PHE
25	d3	100	ASP
25	d3	103	LEU
25	d3	107	PHE
26	d4	6	THR
26	d4	26	ASP
26	d4	34	ASN
26	d4	36	SER
26	d4	40	LEU
26	d4	49	LYS
26	d4	51	GLU
26	d4	58	PHE
26	d4	61	ARG
26	d4	62	THR
26	d4	74	LEU
26	d4	78	SER

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Mol	Chain	Res	Type
26	d4	83	LYS
26	d4	125	LEU
26	d4	133	ASN
27	d5	41	ILE
27	d5	52	LYS
27	d5	53	GLU
27	d5	57	TYR
27	d5	60	VAL
27	d5	77	ARG
27	d5	81	ARG
27	d5	88	ILE
27	d5	92	ILE
27	d5	103	ARG
28	d6	10	ARG
28	d6	11	ASN
28	d6	12	LYS
28	d6	15	ARG
28	d6	24	VAL
28	d6	30	ILE
28	d6	33	ASP
28	d6	39	MET
28	d6	74	CYS
28	d6	82	ARG
28	d6	86	VAL
28	d6	88	SER
28	d6	89	ARG
29	d7	3	LEU
29	d7	21	LEU
29	d7	34	ASP
29	d7	41	LEU
29	d7	42	ASN
29	d7	43	ILE
29	d7	77	THR
29	d7	82	LYS
30	d8	15	VAL
30	d8	18	ARG
30	d8	33	LEU
30	d8	36	THR
30	d8	39	THR
30	d8	49	ARG
30	d8	54	LEU
30	d8	58	GLU

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Mol	Chain	Res	Type
30	d8	61	ARG
30	d8	64	ARG
31	d9	10	HIS
31	d9	12	ARG
31	d9	19	ARG
31	d9	21	CYS
31	d9	28	THR
31	d9	30	LEU
31	d9	32	ARG
31	d9	36	LEU
31	d9	54	LYS
31	d9	56	ARG
32	e0	15	LYS
32	e0	16	SER
32	e0	21	VAL
32	e0	22	GLU
32	e0	24	THR
32	e0	26	LYS
32	e0	29	LYS
32	e0	41	THR
32	e0	43	ARG
32	e0	44	PHE
32	e0	48	THR
32	e0	49	LEU
32	e0	54	ARG
32	e0	56	MET
33	e1	86	THR
33	e1	93	HIS
33	e1	98	VAL
33	e1	99	LYS
33	e1	100	LEU
33	e1	106	TYR
33	e1	113	LYS
33	e1	134	ASN
33	e1	135	HIS
33	e1	147	VAL
33	e1	148	TYR
34	sR	21	THR
34	sR	29	GLN
34	sR	42	LEU
34	sR	58	VAL
34	sR	66	HIS

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Mol	Chain	Res	Type
34	sR	74	THR
34	sR	76	ASP
34	sR	86	ASP
34	sR	106	HIS
34	sR	122	ILE
34	sR	145	LEU
34	sR	167	VAL
34	sR	245	PHE
34	sR	272	ASP
34	sR	286	GLU
34	sR	297	ASP
34	sR	317	THR
35	sM	23	LYS
35	sM	33	LYS
35	sM	43	ASP
35	sM	49	LYS
35	sM	50	ASN
35	sM	68	ARG
35	sM	74	LYS
35	sM	82	THR
39	l2	10	LYS
39	l2	19	HIS
39	l2	32	LEU
39	l2	44	ILE
39	l2	46	LYS
39	l2	47	GLN
39	l2	62	VAL
39	l2	74	GLU
39	l2	80	GLU
39	l2	82	VAL
39	l2	96	LEU
39	l2	101	VAL
39	l2	104	LEU
39	l2	109	GLU
39	l2	114	SER
39	l2	134	VAL
39	l2	137	ILE
39	l2	147	ARG
39	l2	161	ASP
39	l2	165	VAL
39	l2	168	VAL
39	l2	179	LEU

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Mol	Chain	Res	Type
39	12	180	LEU
39	12	193	ARG
39	12	199	THR
39	12	202	VAL
39	12	204	MET
39	12	207	VAL
39	12	241	ARG
39	12	242	ARG
39	12	246	LEU
40	13	3	HIS
40	13	4	ARG
40	13	5	LYS
40	13	7	GLU
40	13	17	LEU
40	13	19	ARG
40	13	24	SER
40	13	30	LYS
40	13	37	ARG
40	13	47	LEU
40	13	50	LYS
40	13	56	ILE
40	13	70	ARG
40	13	79	VAL
40	13	85	VAL
40	13	93	VAL
40	13	103	THR
40	13	104	THR
40	13	110	LEU
40	13	111	SER
40	13	112	ASP
40	13	114	VAL
40	13	116	ARG
40	13	123	TYR
40	13	125	SER
40	13	139	GLN
40	13	140	ASP
40	13	146	ARG
40	13	148	LEU
40	13	156	SER
40	13	167	ARG
40	13	183	LEU
40	13	187	SER

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Mol	Chain	Res	Type
40	l3	188	ILE
40	l3	192	VAL
40	l3	196	ARG
40	l3	208	VAL
40	l3	211	GLN
40	l3	231	HIS
40	l3	232	ARG
40	l3	235	THR
40	l3	249	VAL
40	l3	264	VAL
40	l3	266	ARG
40	l3	274	SER
40	l3	284	ARG
40	l3	296	THR
40	l3	297	SER
40	l3	304	THR
40	l3	324	VAL
40	l3	328	ILE
40	l3	332	ARG
40	l3	335	ILE
40	l3	340	LYS
40	l3	341	SER
40	l3	346	THR
40	l3	348	ARG
40	l3	355	SER
41	l4	25	VAL
41	l4	33	ASP
41	l4	34	ILE
41	l4	54	GLU
41	l4	67	THR
41	l4	69	ARG
41	l4	73	ARG
41	l4	76	ARG
41	l4	92	ASN
41	l4	93	MET
41	l4	120	TYR
41	l4	138	ARG
41	l4	144	LYS
41	l4	150	LEU
41	l4	156	LEU
41	l4	158	SER
41	l4	177	ASP

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Mol	Chain	Res	Type
41	14	179	LEU
41	14	186	LYS
41	14	187	LEU
41	14	191	LYS
41	14	193	LYS
41	14	206	LEU
41	14	220	ARG
41	14	222	VAL
41	14	230	VAL
41	14	246	ARG
41	14	258	LEU
41	14	259	ASP
41	14	272	VAL
41	14	276	LEU
41	14	279	HIS
41	14	301	PRO
41	14	313	LEU
41	14	316	ASN
41	14	319	LYS
41	14	327	LEU
41	14	333	VAL
41	14	347	THR
41	14	361	HIS
42	15	14	SER
42	15	32	GLN
42	15	35	ARG
42	15	51	LEU
42	15	64	ILE
42	15	65	ILE
42	15	93	THR
42	15	109	THR
42	15	110	LEU
42	15	111	GLN
42	15	112	LYS
42	15	118	THR
42	15	120	LYS
42	15	122	VAL
42	15	124	GLU
42	15	133	GLU
42	15	137	ASP
42	15	140	ARG
42	15	144	VAL

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Mol	Chain	Res	Type
42	15	146	LEU
42	15	148	ILE
42	15	155	THR
42	15	185	PHE
42	15	188	GLU
42	15	189	GLU
42	15	194	LEU
42	15	211	LEU
42	15	216	GLU
42	15	218	ARG
42	15	227	LEU
42	15	241	THR
42	15	254	LYS
42	15	258	LYS
42	15	259	LYS
42	15	260	PHE
42	15	263	GLU
42	15	268	GLU
42	15	270	LYS
42	15	275	THR
42	15	278	SER
42	15	279	LYS
42	15	281	GLU
42	15	293	LEU
43	16	13	GLU
43	16	20	LYS
43	16	46	ARG
43	16	48	ARG
43	16	52	VAL
43	16	57	HIS
43	16	62	THR
43	16	64	LEU
43	16	76	LEU
43	16	84	VAL
43	16	89	THR
43	16	92	SER
43	16	131	LYS
43	16	136	GLU
43	16	152	THR
43	16	155	LEU
43	16	157	GLN
43	16	162	SER

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Mol	Chain	Res	Type
43	16	166	LYS
43	16	175	LYS
44	17	22	THR
44	17	38	LYS
44	17	41	ARG
44	17	56	GLU
44	17	60	ARG
44	17	83	LEU
44	17	88	ARG
44	17	93	ASN
44	17	107	ARG
44	17	124	LEU
44	17	130	ILE
44	17	156	ILE
44	17	158	LYS
44	17	159	GLN
44	17	173	LEU
44	17	175	LYS
44	17	179	LEU
44	17	180	SER
44	17	184	LEU
44	17	189	ILE
44	17	199	ASN
44	17	229	PHE
44	17	239	LEU
45	18	43	LYS
45	18	44	ARG
45	18	50	VAL
45	18	57	ARG
45	18	59	GLN
45	18	65	LEU
45	18	71	VAL
45	18	74	THR
45	18	79	GLN
45	18	90	THR
45	18	109	LEU
45	18	132	VAL
45	18	134	TYR
45	18	136	LEU
45	18	160	ILE
45	18	163	VAL
45	18	169	LEU

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Mol	Chain	Res	Type
45	18	172	LYS
45	18	173	MET
45	18	200	LEU
45	18	204	ARG
45	18	213	LYS
45	18	217	THR
45	18	232	HIS
46	19	1	MET
46	19	6	THR
46	19	16	VAL
46	19	17	THR
46	19	18	VAL
46	19	44	THR
46	19	46	THR
46	19	55	VAL
46	19	68	LEU
46	19	70	THR
46	19	80	THR
46	19	118	LEU
46	19	120	ASP
46	19	123	ILE
46	19	130	ASP
46	19	132	VAL
46	19	133	THR
46	19	134	ILE
46	19	144	ILE
46	19	151	VAL
46	19	157	ASN
46	19	162	GLN
46	19	166	ARG
46	19	167	VAL
46	19	173	ARG
46	19	187	ILE
46	19	190	ASP
46	19	191	LEU
47	m0	3	ARG
47	m0	24	ARG
47	m0	28	ASP
47	m0	36	LEU
47	m0	42	THR
47	m0	44	ASP
47	m0	48	LEU

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Mol	Chain	Res	Type
47	m0	52	LEU
47	m0	58	GLU
47	m0	61	SER
47	m0	71	CYS
47	m0	76	MET
47	m0	87	LEU
47	m0	90	ARG
47	m0	99	ILE
47	m0	103	LEU
47	m0	130	ASP
47	m0	137	SER
47	m0	141	LYS
47	m0	156	ARG
47	m0	162	GLN
47	m0	163	GLN
47	m0	167	LEU
47	m0	169	LYS
47	m0	177	ASP
47	m0	183	LYS
47	m0	191	LYS
47	m0	197	VAL
47	m0	205	SER
47	m0	208	ASN
48	m1	6	GLN
48	m1	10	ARG
48	m1	12	LEU
48	m1	13	LYS
48	m1	16	LYS
48	m1	23	VAL
48	m1	30	LEU
48	m1	46	VAL
48	m1	92	ARG
48	m1	99	THR
48	m1	101	ASN
48	m1	107	ASP
48	m1	111	ASP
48	m1	112	LEU
48	m1	129	VAL
48	m1	137	ARG
48	m1	139	THR
48	m1	147	THR
48	m1	148	VAL

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Mol	Chain	Res	Type
48	m1	158	ASP
48	m1	159	THR
48	m1	165	GLN
49	m3	8	PRO
49	m3	13	HIS
49	m3	52	ASP
49	m3	54	LEU
49	m3	55	ARG
49	m3	58	VAL
49	m3	67	ARG
49	m3	69	VAL
49	m3	73	ARG
49	m3	86	THR
49	m3	97	VAL
49	m3	100	ARG
49	m3	103	ASN
49	m3	118	GLU
49	m3	128	ARG
49	m3	131	LYS
49	m3	147	ILE
49	m3	149	GLN
49	m3	154	VAL
49	m3	157	ARG
49	m3	164	GLU
49	m3	168	ARG
49	m3	170	LEU
49	m3	176	GLU
49	m3	180	ARG
49	m3	189	GLU
49	m3	194	GLU
50	m4	4	ASP
50	m4	8	LYS
50	m4	11	ASN
50	m4	15	VAL
50	m4	28	SER
50	m4	42	LYS
50	m4	53	VAL
50	m4	69	THR
50	m4	72	LEU
50	m4	80	THR
50	m4	106	ARG
50	m4	107	GLU

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Mol	Chain	Res	Type
50	m4	108	ARG
50	m4	113	THR
51	m5	5	LYS
51	m5	7	LEU
51	m5	10	LEU
51	m5	12	ARG
51	m5	17	ASP
51	m5	19	LEU
51	m5	22	LEU
51	m5	27	VAL
51	m5	49	ARG
51	m5	71	ARG
51	m5	80	THR
51	m5	86	ASN
51	m5	106	VAL
51	m5	109	ARG
51	m5	117	ASN
51	m5	138	GLN
51	m5	142	ILE
51	m5	153	ASP
51	m5	176	LYS
51	m5	182	ASN
51	m5	183	THR
52	m6	4	GLU
52	m6	22	VAL
52	m6	34	VAL
52	m6	46	GLU
52	m6	60	LYS
52	m6	74	ARG
52	m6	85	ARG
52	m6	106	GLU
52	m6	116	LYS
52	m6	118	VAL
52	m6	124	LEU
52	m6	126	VAL
52	m6	129	LEU
52	m6	134	LYS
52	m6	143	THR
52	m6	151	ASP
52	m6	160	ARG
52	m6	162	VAL
52	m6	175	THR

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Mol	Chain	Res	Type
52	m6	178	VAL
52	m6	197	LEU
53	m7	7	THR
53	m7	9	THR
53	m7	24	VAL
53	m7	32	THR
53	m7	41	LEU
53	m7	47	TYR
53	m7	52	LEU
53	m7	79	THR
53	m7	80	LYS
53	m7	89	LYS
53	m7	107	LEU
53	m7	112	LEU
53	m7	119	VAL
53	m7	126	ARG
53	m7	135	ARG
53	m7	136	ILE
53	m7	147	GLU
53	m7	148	LEU
53	m7	150	VAL
54	m8	3	ILE
54	m8	7	SER
54	m8	12	ARG
54	m8	17	THR
54	m8	22	ASP
54	m8	26	LEU
54	m8	31	LYS
54	m8	32	LEU
54	m8	34	THR
54	m8	57	ILE
54	m8	66	ARG
54	m8	69	ARG
54	m8	81	VAL
54	m8	86	THR
54	m8	98	LYS
54	m8	127	LEU
54	m8	135	GLN
54	m8	138	LEU
54	m8	161	LYS
54	m8	165	ILE
54	m8	170	ARG

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Mol	Chain	Res	Type
54	m8	174	ARG
54	m8	178	ARG
54	m8	181	SER
55	m9	7	GLN
55	m9	10	LEU
55	m9	21	LYS
55	m9	29	THR
55	m9	30	SER
55	m9	34	GLN
55	m9	43	LYS
55	m9	49	THR
55	m9	55	VAL
55	m9	63	THR
55	m9	74	ARG
55	m9	88	ARG
55	m9	138	LEU
55	m9	152	GLU
55	m9	164	LEU
55	m9	166	ASN
55	m9	173	ARG
55	m9	180	LYS
55	m9	182	ASP
55	m9	186	LYS
56	n0	1	MET
56	n0	3	HIS
56	n0	8	GLN
56	n0	21	GLU
56	n0	73	LYS
56	n0	80	ARG
56	n0	87	THR
56	n0	96	ASP
56	n0	97	VAL
56	n0	100	VAL
56	n0	105	THR
56	n0	107	TYR
56	n0	130	GLU
56	n0	132	THR
56	n0	136	LYS
56	n0	137	ARG
56	n0	145	THR
56	n0	148	LEU
56	n0	155	ARG

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Mol	Chain	Res	Type
56	n0	160	THR
56	n0	161	LYS
56	n0	162	THR
56	n0	172	TYR
57	n1	25	VAL
57	n1	27	LEU
57	n1	33	VAL
57	n1	47	SER
57	n1	71	SER
57	n1	78	LYS
57	n1	80	VAL
57	n1	83	ARG
57	n1	89	LEU
57	n1	96	ILE
57	n1	103	GLN
57	n1	126	VAL
57	n1	128	LEU
57	n1	130	ARG
57	n1	131	GLN
57	n1	138	SER
57	n1	139	ARG
57	n1	141	VAL
57	n1	143	THR
57	n1	150	THR
57	n1	157	GLU
57	n1	158	THR
57	n1	159	PHE
58	n2	27	VAL
58	n2	29	ASP
58	n2	43	VAL
58	n2	57	THR
58	n2	58	GLU
58	n2	68	THR
58	n2	74	LYS
58	n2	75	TYR
58	n2	90	ARG
58	n2	94	ARG
58	n2	100	THR
58	n2	108	TYR
59	n3	4	ASN
59	n3	7	GLN
59	n3	12	ARG

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Mol	Chain	Res	Type
59	n3	13	ILE
59	n3	14	SER
59	n3	48	ARG
59	n3	68	GLU
59	n3	73	VAL
59	n3	74	MET
59	n3	84	SER
59	n3	93	LEU
59	n3	96	GLU
59	n3	109	MET
60	n4	5	ILE
60	n4	39	LEU
60	n4	42	GLN
60	n4	47	ARG
60	n4	54	LEU
61	n5	24	LEU
61	n5	27	ARG
61	n5	38	LEU
61	n5	39	LYS
61	n5	56	ARG
61	n5	63	ILE
61	n5	71	THR
61	n5	108	LEU
61	n5	115	ARG
61	n5	125	ARG
61	n5	134	ASP
61	n5	135	ILE
61	n5	137	ASN
61	n5	138	ARG
62	n6	4	GLN
62	n6	12	ARG
62	n6	13	ARG
62	n6	32	SER
62	n6	37	LYS
62	n6	45	ILE
62	n6	48	LEU
62	n6	50	ILE
62	n6	51	ARG
62	n6	56	VAL
62	n6	59	VAL
62	n6	66	GLN
62	n6	74	TYR

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Mol	Chain	Res	Type
62	n6	83	ASP
62	n6	99	LEU
62	n6	115	ARG
63	n7	14	VAL
63	n7	17	ARG
63	n7	24	VAL
63	n7	36	HIS
63	n7	46	ILE
63	n7	52	LYS
63	n7	56	LYS
63	n7	57	HIS
63	n7	72	ILE
63	n7	81	LEU
63	n7	84	ARG
63	n7	90	GLU
63	n7	92	PHE
63	n7	100	THR
63	n7	118	PHE
63	n7	126	LYS
63	n7	128	GLN
63	n7	134	LEU
64	n8	4	ARG
64	n8	6	THR
64	n8	8	THR
64	n8	10	LYS
64	n8	15	VAL
64	n8	19	LYS
64	n8	24	LYS
64	n8	26	ARG
64	n8	27	LYS
64	n8	42	ARG
64	n8	56	VAL
64	n8	60	TYR
64	n8	65	GLN
64	n8	91	LEU
64	n8	123	VAL
64	n8	132	LYS
64	n8	139	ARG
65	n9	6	ASN
65	n9	13	THR
65	n9	18	ARG
65	n9	38	LYS

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Mol	Chain	Res	Type
65	n9	59	LYS
66	o0	12	GLN
66	o0	28	LYS
66	o0	29	SER
66	o0	40	LYS
66	o0	41	LEU
66	o0	50	VAL
66	o0	61	MET
66	o0	63	SER
66	o0	66	LYS
66	o0	86	ARG
66	o0	87	VAL
66	o0	103	THR
67	o1	13	THR
67	o1	16	LEU
67	o1	24	SER
67	o1	26	LYS
67	o1	31	ARG
67	o1	41	LYS
67	o1	42	LEU
67	o1	44	MET
67	o1	50	ARG
67	o1	62	ARG
67	o1	68	GLU
67	o1	83	GLU
67	o1	91	SER
67	o1	93	VAL
67	o1	96	VAL
67	o1	102	LYS
67	o1	106	THR
67	o1	110	GLU
68	o2	3	SER
68	o2	6	HIS
68	o2	19	ARG
68	o2	24	ARG
68	o2	27	ARG
68	o2	33	ARG
68	o2	38	ILE
68	o2	49	ASN
68	o2	51	SER
68	o2	59	SER
68	o2	61	LYS

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Mol	Chain	Res	Type
68	o2	73	THR
68	o2	75	LEU
68	o2	81	ASP
68	o2	82	LEU
68	o2	84	THR
68	o2	91	THR
68	o2	95	GLU
68	o2	103	LYS
68	o2	109	LEU
68	o2	125	ARG
68	o2	126	LEU
69	o3	3	GLU
69	o3	15	SER
69	o3	20	LYS
69	o3	21	ARG
69	o3	28	SER
69	o3	31	LYS
69	o3	37	THR
69	o3	45	LEU
69	o3	74	THR
69	o3	80	VAL
69	o3	81	VAL
69	o3	87	ASN
69	o3	98	VAL
69	o3	105	SER
70	o4	5	VAL
70	o4	19	LYS
70	o4	20	ILE
70	o4	35	VAL
70	o4	58	ARG
70	o4	65	VAL
70	o4	71	THR
70	o4	83	ASN
70	o4	104	VAL
71	o5	11	THR
71	o5	16	GLN
71	o5	21	LEU
71	o5	38	ARG
71	o5	45	LYS
71	o5	46	THR
71	o5	47	VAL
71	o5	49	LYS

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Mol	Chain	Res	Type
71	o5	56	THR
71	o5	69	LEU
71	o5	79	ASP
71	o5	85	THR
71	o5	101	THR
72	o6	7	ILE
72	o6	9	ILE
72	o6	11	LEU
72	o6	15	LYS
72	o6	17	VAL
72	o6	21	THR
72	o6	26	ILE
72	o6	29	LYS
72	o6	35	ASN
72	o6	36	ARG
72	o6	43	LEU
72	o6	45	ARG
72	o6	57	LEU
72	o6	58	ILE
72	o6	60	LEU
72	o6	68	ARG
72	o6	72	VAL
72	o6	76	ARG
72	o6	81	THR
72	o6	94	ILE
72	o6	98	ARG
73	o7	14	LYS
73	o7	25	ARG
73	o7	33	THR
73	o7	46	SER
73	o7	59	THR
73	o7	65	ARG
73	o7	67	LEU
73	o7	75	LYS
74	o8	8	ILE
74	o8	12	LEU
74	o8	14	LEU
74	o8	24	THR
74	o8	41	THR
74	o8	53	THR
74	o8	54	LEU
74	o8	61	LYS

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Mol	Chain	Res	Type
74	o8	64	LYS
74	o8	66	ILE
75	o9	4	GLN
75	o9	21	ARG
75	o9	46	ARG
76	q0	79	GLU
76	q0	85	LEU
76	q0	93	LYS
76	q0	96	CYS
76	q0	99	CYS
76	q0	112	LYS
76	q0	113	ARG
76	q0	127	LEU
77	q1	9	ARG
77	q1	10	THR
77	q1	13	LEU
77	q1	17	ARG
77	q1	18	ARG
77	q1	21	ARG
78	q2	7	THR
78	q2	8	ARG
78	q2	10	THR
78	q2	19	LYS
78	q2	26	THR
78	q2	57	VAL
78	q2	61	LYS
78	q2	66	LYS
78	q2	67	LYS
78	q2	68	VAL
78	q2	74	CYS
78	q2	79	THR
78	q2	82	GLN
78	q2	84	THR
78	q2	85	LEU
78	q2	93	LEU
78	q2	100	LYS
78	q2	105	GLN
79	q3	16	VAL
79	q3	42	CYS
79	q3	46	THR
79	q3	49	ARG
79	q3	54	ILE

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Mol	Chain	Res	Type
79	q3	56	THR
79	q3	58	SER
79	q3	73	THR
79	q3	78	THR
79	q3	89	MET
81	p0	4	ILE
81	p0	5	ARG
81	p0	15	LEU
81	p0	39	HIS
81	p0	48	ARG
81	p0	52	LEU
81	p0	67	LEU
81	p0	70	LEU
81	p0	81	LYS
81	p0	93	LEU
81	p0	97	LYS
81	p0	105	VAL
83	f	37	ARG
83	f	51	LYS
83	f	52	HIS
83	f	54	HIS
83	f	58	HIS
83	f	78	HIS
83	f	79	ASN
83	f	80	MET
83	f	87	ARG
83	f	102	LEU
83	f	103	MET
83	f	105	MET
83	f	106	ASP
83	f	109	THR
83	f	112	ASP
83	f	120	LEU
83	f	138	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
2	S0	163	ASN
3	S1	209	ASN
4	S2	77	GLN
20	C8	99	HIS

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Mol	Chain	Res	Type
23	D1	3	ASN
25	D3	65	ASN
25	D3	79	ASN
26	D4	106	GLN
34	SR	248	ASN
41	L4	311	HIS
42	L5	40	HIS
46	L9	49	ASN
46	L9	116	ASN
47	M0	144	ASN
48	M1	95	ASN
51	M5	178	HIS
51	M5	194	GLN
54	M8	73	GLN
57	N1	26	HIS
3	s1	157	GLN
4	s2	94	GLN
7	s5	34	GLN
9	s7	71	HIS
9	s7	74	GLN
12	c0	32	HIS
15	c3	69	ASN
18	c6	139	GLN
24	d2	15	ASN
26	d4	34	ASN
32	e0	17	GLN
33	e1	93	HIS
34	sR	268	GLN
45	l8	59	GLN
51	m5	87	GLN
52	m6	72	HIS
53	m7	137	ASN
64	n8	25	HIS
65	n9	12	GLN
75	o9	25	GLN
78	q2	82	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	2	1776/1800 (98%)	455 (25%)	40 (2%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	6	1791/1800 (99%)	447 (24%)	29 (1%)
36	1	3145/3396 (92%)	672 (21%)	53 (1%)
36	5	3163/3396 (93%)	649 (20%)	52 (1%)
37	3	120/121 (99%)	11 (9%)	0
37	7	120/121 (99%)	13 (10%)	0
38	4	157/158 (99%)	32 (20%)	3 (1%)
38	8	157/158 (99%)	34 (21%)	0
All	All	10429/10950 (95%)	2313 (22%)	177 (1%)

All (2313) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	2	2	A
1	2	4	C
1	2	6	G
1	2	17	C
1	2	23	G
1	2	25	C
1	2	26	A
1	2	27	U
1	2	34	G
1	2	42	G
1	2	45	U
1	2	46	A
1	2	47	A
1	2	57	G
1	2	60	U
1	2	67	A
1	2	68	A
1	2	69	G
1	2	72	A
1	2	73	U
1	2	74	U
1	2	75	U
1	2	77	U
1	2	95	G
1	2	96	G
1	2	104	A
1	2	111	U
1	2	114	C
1	2	121	U
1	2	127	G

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Mol	Chain	Res	Type
1	2	128	U
1	2	131	C
1	2	132	U
1	2	133	U
1	2	134	U
1	2	135	A
1	2	136	C
1	2	137	U
1	2	141	U
1	2	144	U
1	2	145	A
1	2	146	U
1	2	153	G
1	2	159	U
1	2	160	C
1	2	176	C
1	2	178	U
1	2	179	A
1	2	184	C
1	2	185	U
1	2	187	G
1	2	190	C
1	2	191	C
1	2	192	U
1	2	194	U
1	2	195	G
1	2	197	A
1	2	200	A
1	2	215	A
1	2	217	A
1	2	218	A
1	2	219	A
1	2	221	A
1	2	226	A
1	2	227	U
1	2	228	G
1	2	231	U
1	2	233	C
1	2	234	G
1	2	235	G
1	2	236	A
1	2	239	C

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Mol	Chain	Res	Type
1	2	240	U
1	2	241	U
1	2	242	U
1	2	245	U
1	2	249	U
1	2	250	C
1	2	257	A
1	2	261	U
1	2	262	U
1	2	265	A
1	2	271	A
1	2	272	U
1	2	273	G
1	2	274	G
1	2	275	C
1	2	276	C
1	2	277	U
1	2	278	U
1	2	279	G
1	2	280	U
1	2	281	G
1	2	288	A
1	2	290	G
1	2	299	A
1	2	302	U
1	2	309	C
1	2	314	C
1	2	316	A
1	2	319	U
1	2	321	C
1	2	329	G
1	2	333	A
1	2	337	G
1	2	338	C
1	2	352	A
1	2	358	U
1	2	359	A
1	2	360	A
1	2	361	C
1	2	380	U
1	2	390	G
1	2	400	A

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Mol	Chain	Res	Type
1	2	402	C
1	2	404	G
1	2	416	A
1	2	418	G
1	2	419	G
1	2	421	A
1	2	424	C
1	2	425	A
1	2	426	G
1	2	428	A
1	2	434	G
1	2	437	A
1	2	439	U
1	2	440	U
1	2	444	C
1	2	445	A
1	2	447	U
1	2	448	C
1	2	455	C
1	2	468	A
1	2	469	C
1	2	477	A
1	2	483	A
1	2	484	C
1	2	485	A
1	2	486	G
1	2	488	G
1	2	493	U
1	2	494	U
1	2	495	C
1	2	496	G
1	2	497	G
1	2	498	G
1	2	499	U
1	2	500	C
1	2	502	U
1	2	503	G
1	2	504	U
1	2	505	A
1	2	506	A
1	2	507	U
1	2	508	U

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Mol	Chain	Res	Type
1	2	510	G
1	2	511	A
1	2	513	U
1	2	514	G
1	2	515	A
1	2	516	G
1	2	518	A
1	2	520	A
1	2	527	A
1	2	529	A
1	2	532	U
1	2	534	A
1	2	538	A
1	2	539	G
1	2	540	G
1	2	541	A
1	2	542	A
1	2	543	C
1	2	544	A
1	2	556	A
1	2	557	G
1	2	558	U
1	2	559	C
1	2	565	C
1	2	574	G
1	2	579	A
1	2	580	A
1	2	585	A
1	2	594	A
1	2	595	G
1	2	606	A
1	2	611	U
1	2	619	A
1	2	620	A
1	2	621	A
1	2	622	A
1	2	623	A
1	2	624	G
1	2	628	G
1	2	639	U
1	2	640	U
1	2	648	G

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Mol	Chain	Res	Type
1	2	650	U
1	2	653	C
1	2	655	G
1	2	656	G
1	2	657	U
1	2	658	C
1	2	680	U
1	2	684	A
1	2	686	C
1	2	687	G
1	2	691	C
1	2	694	U
1	2	696	C
1	2	697	C
1	2	700	C
1	2	701	U
1	2	702	G
1	2	703	G
1	2	704	C
1	2	705	U
1	2	706	A
1	2	707	A
1	2	709	C
1	2	710	U
1	2	712	G
1	2	714	G
1	2	717	C
1	2	718	U
1	2	719	U
1	2	721	U
1	2	722	G
1	2	723	G
1	2	725	U
1	2	727	U
1	2	728	U
1	2	731	C
1	2	732	G
1	2	734	A
1	2	735	C
1	2	736	C
1	2	737	A
1	2	738	G

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Mol	Chain	Res	Type
1	2	742	U
1	2	743	U
1	2	754	A
1	2	755	A
1	2	756	A
1	2	765	G
1	2	766	U
1	2	774	A
1	2	775	G
1	2	778	G
1	2	781	U
1	2	783	G
1	2	784	C
1	2	789	A
1	2	794	U
1	2	795	U
1	2	812	A
1	2	815	G
1	2	816	G
1	2	818	C
1	2	819	G
1	2	820	U
1	2	821	U
1	2	824	G
1	2	829	A
1	2	830	U
1	2	831	U
1	2	833	U
1	2	839	U
1	2	846	G
1	2	856	A
1	2	860	U
1	2	863	A
1	2	876	G
1	2	886	U
1	2	898	A
1	2	912	U
1	2	914	G
1	2	933	A
1	2	935	U
1	2	940	A
1	2	942	G

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Mol	Chain	Res	Type
1	2	944	A
1	2	951	A
1	2	954	G
1	2	960	U
1	2	966	A
1	2	969	C
1	2	992	A
1	2	993	A
1	2	995	A
1	2	1003	A
1	2	1004	U
1	2	1005	A
1	2	1022	C
1	2	1026	A
1	2	1028	C
1	2	1039	A
1	2	1040	G
1	2	1052	U
1	2	1053	G
1	2	1056	U
1	2	1057	U
1	2	1058	U
1	2	1060	U
1	2	1061	A
1	2	1074	G
1	2	1080	U
1	2	1082	C
1	2	1091	A
1	2	1092	A
1	2	1096	C
1	2	1097	U
1	2	1098	U
1	2	1100	G
1	2	1109	G
1	2	1111	G
1	2	1138	A
1	2	1146	G
1	2	1150	G
1	2	1151	A
1	2	1155	G
1	2	1158	C
1	2	1160	A

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Mol	Chain	Res	Type
1	2	1167	G
1	2	1185	U
1	2	1194	A
1	2	1196	A
1	2	1197	C
1	2	1199	G
1	2	1200	G
1	2	1208	A
1	2	1217	A
1	2	1218	G
1	2	1227	A
1	2	1229	G
1	2	1243	G
1	2	1244	A
1	2	1245	G
1	2	1250	U
1	2	1251	U
1	2	1258	U
1	2	1284	C
1	2	1286	U
1	2	1301	U
1	2	1311	U
1	2	1314	U
1	2	1315	U
1	2	1321	A
1	2	1334	U
1	2	1337	A
1	2	1340	U
1	2	1341	A
1	2	1344	A
1	2	1345	A
1	2	1346	A
1	2	1347	U
1	2	1348	A
1	2	1354	G
1	2	1355	C
1	2	1361	U
1	2	1363	U
1	2	1364	G
1	2	1370	U
1	2	1371	A
1	2	1372	U

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Mol	Chain	Res	Type
1	2	1390	U
1	2	1398	U
1	2	1399	C
1	2	1413	U
1	2	1414	U
1	2	1415	U
1	2	1427	A
1	2	1428	G
1	2	1432	U
1	2	1436	A
1	2	1437	U
1	2	1444	A
1	2	1446	A
1	2	1449	U
1	2	1456	C
1	2	1458	G
1	2	1459	C
1	2	1460	A
1	2	1461	C
1	2	1471	A
1	2	1473	U
1	2	1474	G
1	2	1477	G
1	2	1481	C
1	2	1482	C
1	2	1486	G
1	2	1489	U
1	2	1490	C
1	2	1491	U
1	2	1492	A
1	2	1493	A
1	2	1499	G
1	2	1506	G
1	2	1515	A
1	2	1516	A
1	2	1517	U
1	2	1521	G
1	2	1523	G
1	2	1524	A
1	2	1535	U
1	2	1536	G
1	2	1537	C

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Mol	Chain	Res	Type
1	2	1538	U
1	2	1542	G
1	2	1557	U
1	2	1559	A
1	2	1569	A
1	2	1574	G
1	2	1582	U
1	2	1584	G
1	2	1601	G
1	2	1616	G
1	2	1619	C
1	2	1624	C
1	2	1626	U
1	2	1631	A
1	2	1635	A
1	2	1636	C
1	2	1657	U
1	2	1658	G
1	2	1680	G
1	2	1683	C
1	2	1684	U
1	2	1685	G
1	2	1689	A
1	2	1698	G
1	2	1699	G
1	2	1700	C
1	2	1701	A
1	2	1702	A
1	2	1703	C
1	2	1712	A
1	2	1713	G
1	2	1714	A
1	2	1731	A
1	2	1760	G
1	2	1762	A
1	2	1769	U
1	2	1770	U
1	2	1771	U
1	2	1780	G
1	2	1782	A
1	2	1783	C
1	2	1792	G

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Mol	Chain	Res	Type
1	2	1793	G
1	2	1794	A
1	2	1795	U
1	2	1796	C
1	2	1797	A
36	1	15	C
36	1	16	A
36	1	21	G
36	1	24	G
36	1	26	A
36	1	40	A
36	1	43	A
36	1	49	A
36	1	59	G
36	1	60	A
36	1	65	A
36	1	66	A
36	1	67	A
36	1	68	C
36	1	74	G
36	1	83	U
36	1	92	G
36	1	93	C
36	1	99	A
36	1	108	A
36	1	109	A
36	1	110	G
36	1	111	C
36	1	113	C
36	1	116	A
36	1	117	U
36	1	121	A
36	1	122	A
36	1	131	C
36	1	133	U
36	1	136	G
36	1	140	C
36	1	154	U
36	1	155	G
36	1	156	G
36	1	157	A
36	1	173	G

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Mol	Chain	Res	Type
36	1	175	C
36	1	187	A
36	1	190	U
36	1	191	U
36	1	201	A
36	1	206	G
36	1	210	U
36	1	213	A
36	1	218	G
36	1	219	A
36	1	220	G
36	1	239	G
36	1	240	U
36	1	241	G
36	1	243	G
36	1	245	U
36	1	249	U
36	1	250	U
36	1	252	U
36	1	253	A
36	1	256	G
36	1	269	G
36	1	282	G
36	1	283	G
36	1	286	U
36	1	287	G
36	1	288	C
36	1	295	A
36	1	298	U
36	1	299	G
36	1	305	U
36	1	315	C
36	1	323	A
36	1	329	U
36	1	330	G
36	1	338	A
36	1	339	C
36	1	349	A
36	1	350	C
36	1	366	A
36	1	376	G
36	1	380	U

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Mol	Chain	Res	Type
36	1	398	A
36	1	399	A
36	1	401	U
36	1	402	A
36	1	403	C
36	1	414	U
36	1	421	G
36	1	422	A
36	1	439	C
36	1	440	A
36	1	495	G
36	1	498	A
36	1	520	U
36	1	521	A
36	1	523	A
36	1	535	G
36	1	543	C
36	1	544	C
36	1	546	C
36	1	547	G
36	1	548	G
36	1	550	A
36	1	552	G
36	1	555	U
36	1	556	U
36	1	557	A
36	1	559	A
36	1	568	G
36	1	578	A
36	1	579	G
36	1	588	G
36	1	592	A
36	1	602	A
36	1	604	G
36	1	609	G
36	1	611	A
36	1	619	A
36	1	620	U
36	1	621	A
36	1	636	C
36	1	649	A
36	1	660	A

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Mol	Chain	Res	Type
36	1	677	A
36	1	681	U
36	1	682	U
36	1	689	U
36	1	690	A
36	1	691	A
36	1	705	A
36	1	708	G
36	1	709	A
36	1	712	G
36	1	715	A
36	1	716	A
36	1	718	G
36	1	764	U
36	1	765	C
36	1	766	U
36	1	767	U
36	1	768	C
36	1	776	U
36	1	777	U
36	1	781	G
36	1	785	G
36	1	792	G
36	1	806	A
36	1	817	A
36	1	830	A
36	1	837	A
36	1	849	C
36	1	861	C
36	1	873	C
36	1	874	U
36	1	879	U
36	1	890	C
36	1	896	A
36	1	907	G
36	1	908	G
36	1	914	A
36	1	916	G
36	1	917	A
36	1	921	A
36	1	923	C
36	1	924	G

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Mol	Chain	Res	Type
36	1	937	G
36	1	943	U
36	1	944	C
36	1	953	G
36	1	959	C
36	1	960	U
36	1	962	A
36	1	963	G
36	1	964	G
36	1	974	G
36	1	979	U
36	1	980	A
36	1	981	U
36	1	982	C
36	1	993	G
36	1	994	G
36	1	1001	G
36	1	1002	A
36	1	1010	G
36	1	1013	G
36	1	1016	C
36	1	1017	C
36	1	1018	G
36	1	1020	G
36	1	1024	G
36	1	1025	A
36	1	1029	G
36	1	1037	C
36	1	1038	C
36	1	1047	A
36	1	1049	C
36	1	1063	G
36	1	1064	A
36	1	1065	A
36	1	1071	U
36	1	1072	G
36	1	1081	U
36	1	1082	U
36	1	1093	A
36	1	1094	U
36	1	1095	U
36	1	1097	G

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Mol	Chain	Res	Type
36	1	1098	A
36	1	1103	A
36	1	1104	G
36	1	1117	G
36	1	1128	U
36	1	1131	G
36	1	1144	U
36	1	1153	A
36	1	1156	C
36	1	1159	A
36	1	1178	G
36	1	1179	A
36	1	1180	A
36	1	1181	U
36	1	1182	A
36	1	1190	A
36	1	1191	U
36	1	1192	C
36	1	1201	C
36	1	1206	G
36	1	1209	G
36	1	1212	A
36	1	1217	A
36	1	1222	G
36	1	1227	C
36	1	1232	C
36	1	1235	U
36	1	1236	G
36	1	1237	G
36	1	1240	A
36	1	1241	U
36	1	1243	G
36	1	1245	A
36	1	1246	G
36	1	1248	C
36	1	1249	G
36	1	1253	U
36	1	1254	C
36	1	1258	U
36	1	1262	G
36	1	1263	A
36	1	1264	G

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Mol	Chain	Res	Type
36	1	1266	G
36	1	1270	A
36	1	1271	A
36	1	1274	A
36	1	1278	A
36	1	1279	C
36	1	1285	G
36	1	1287	A
36	1	1305	U
36	1	1308	A
36	1	1309	U
36	1	1318	A
36	1	1330	A
36	1	1331	U
36	1	1332	A
36	1	1345	G
36	1	1348	U
36	1	1349	G
36	1	1350	A
36	1	1351	U
36	1	1352	A
36	1	1353	U
36	1	1354	G
36	1	1355	A
36	1	1356	U
36	1	1357	G
36	1	1386	A
36	1	1387	G
36	1	1392	G
36	1	1398	U
36	1	1399	A
36	1	1400	G
36	1	1408	G
36	1	1417	G
36	1	1418	A
36	1	1419	A
36	1	1421	G
36	1	1429	G
36	1	1431	G
36	1	1433	A
36	1	1434	G
36	1	1437	C

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Mol	Chain	Res	Type
36	1	1446	A
36	1	1481	A
36	1	1482	A
36	1	1483	G
36	1	1485	G
36	1	1502	C
36	1	1508	C
36	1	1525	G
36	1	1527	C
36	1	1535	A
36	1	1556	C
36	1	1560	G
36	1	1562	C
36	1	1563	C
36	1	1567	U
36	1	1568	U
36	1	1569	U
36	1	1570	U
36	1	1571	A
36	1	1576	G
36	1	1579	C
36	1	1580	A
36	1	1582	C
36	1	1583	A
36	1	1587	A
36	1	1589	A
36	1	1592	G
36	1	1605	A
36	1	1620	U
36	1	1621	A
36	1	1629	U
36	1	1633	C
36	1	1639	C
36	1	1641	U
36	1	1642	A
36	1	1643	A
36	1	1644	C
36	1	1645	U
36	1	1655	G
36	1	1657	C
36	1	1658	G
36	1	1675	G

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Mol	Chain	Res	Type
36	1	1683	A
36	1	1688	U
36	1	1702	U
36	1	1713	G
36	1	1715	A
36	1	1716	U
36	1	1717	U
36	1	1724	U
36	1	1729	A
36	1	1736	G
36	1	1740	U
36	1	1742	U
36	1	1750	A
36	1	1751	G
36	1	1753	G
36	1	1761	C
36	1	1762	C
36	1	1763	U
36	1	1764	U
36	1	1765	U
36	1	1766	G
36	1	1767	C
36	1	1770	G
36	1	1775	G
36	1	1778	G
36	1	1780	G
36	1	1781	C
36	1	1797	A
36	1	1798	A
36	1	1810	A
36	1	1814	A
36	1	1816	A
36	1	1819	U
36	1	1820	U
36	1	1821	U
36	1	1834	U
36	1	1839	A
36	1	1841	A
36	1	1842	A
36	1	1846	C
36	1	1849	C
36	1	1850	A

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Mol	Chain	Res	Type
36	1	1871	U
36	1	1879	A
36	1	1880	U
36	1	1881	A
36	1	1906	G
36	1	1937	U
36	1	1948	G
36	1	1952	G
36	1	1953	G
36	1	1954	G
36	1	2094	C
36	1	2095	G
36	1	2101	C
36	1	2102	U
36	1	2111	G
36	1	2113	A
36	1	2115	G
36	1	2116	G
36	1	2120	A
36	1	2121	G
36	1	2122	G
36	1	2130	G
36	1	2131	A
36	1	2140	U
36	1	2144	A
36	1	2158	A
36	1	2169	G
36	1	2170	U
36	1	2187	G
36	1	2188	A
36	1	2205	U
36	1	2207	A
36	1	2208	A
36	1	2209	U
36	1	2210	G
36	1	2215	A
36	1	2242	A
36	1	2244	A
36	1	2249	G
36	1	2250	G
36	1	2252	A
36	1	2255	A

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Mol	Chain	Res	Type
36	1	2256	A
36	1	2260	U
36	1	2272	G
36	1	2281	A
36	1	2282	U
36	1	2283	G
36	1	2287	C
36	1	2298	U
36	1	2299	A
36	1	2307	G
36	1	2310	U
36	1	2313	A
36	1	2314	U
36	1	2315	G
36	1	2334	U
36	1	2335	G
36	1	2336	U
36	1	2361	A
36	1	2372	A
36	1	2373	A
36	1	2374	C
36	1	2375	G
36	1	2382	G
36	1	2385	G
36	1	2393	G
36	1	2394	G
36	1	2397	A
36	1	2398	A
36	1	2401	A
36	1	2402	A
36	1	2403	G
36	1	2404	A
36	1	2411	U
36	1	2418	G
36	1	2419	A
36	1	2437	G
36	1	2444	C
36	1	2445	A
36	1	2502	A
36	1	2503	G
36	1	2508	U
36	1	2514	U

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Mol	Chain	Res	Type
36	1	2515	A
36	1	2522	G
36	1	2523	A
36	1	2526	C
36	1	2532	U
36	1	2533	G
36	1	2534	G
36	1	2537	U
36	1	2538	U
36	1	2539	C
36	1	2540	A
36	1	2541	U
36	1	2542	U
36	1	2543	U
36	1	2549	G
36	1	2552	C
36	1	2554	A
36	1	2555	G
36	1	2560	C
36	1	2561	A
36	1	2562	A
36	1	2568	C
36	1	2569	A
36	1	2570	U
36	1	2571	U
36	1	2572	C
36	1	2573	G
36	1	2581	U
36	1	2585	G
36	1	2586	G
36	1	2587	U
36	1	2593	A
36	1	2594	C
36	1	2606	G
36	1	2607	G
36	1	2614	G
36	1	2617	U
36	1	2626	A
36	1	2637	A
36	1	2642	A
36	1	2652	U
36	1	2656	A

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Mol	Chain	Res	Type
36	1	2657	A
36	1	2672	G
36	1	2674	A
36	1	2677	G
36	1	2689	A
36	1	2691	A
36	1	2694	A
36	1	2696	A
36	1	2704	A
36	1	2705	A
36	1	2706	G
36	1	2714	G
36	1	2728	G
36	1	2729	U
36	1	2740	A
36	1	2749	G
36	1	2752	U
36	1	2753	G
36	1	2755	C
36	1	2760	C
36	1	2762	A
36	1	2771	U
36	1	2772	C
36	1	2777	G
36	1	2778	G
36	1	2787	G
36	1	2796	G
36	1	2797	C
36	1	2799	A
36	1	2800	G
36	1	2801	A
36	1	2802	A
36	1	2810	C
36	1	2814	G
36	1	2816	G
36	1	2817	A
36	1	2818	U
36	1	2837	A
36	1	2838	A
36	1	2842	U
36	1	2843	U
36	1	2845	A

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Mol	Chain	Res	Type
36	1	2847	A
36	1	2853	A
36	1	2856	G
36	1	2858	U
36	1	2860	U
36	1	2861	U
36	1	2867	C
36	1	2871	G
36	1	2872	A
36	1	2887	A
36	1	2898	G
36	1	2899	C
36	1	2923	U
36	1	2932	U
36	1	2935	U
36	1	2936	A
36	1	2941	A
36	1	2942	C
36	1	2947	G
36	1	2951	G
36	1	2960	C
36	1	2971	A
36	1	2972	G
36	1	2977	G
36	1	2983	C
36	1	2990	G
36	1	2992	U
36	1	2996	U
36	1	2997	G
36	1	3012	A
36	1	3025	C
36	1	3030	G
36	1	3049	A
36	1	3056	U
36	1	3059	G
36	1	3074	G
36	1	3078	U
36	1	3079	U
36	1	3080	G
36	1	3086	A
36	1	3087	A
36	1	3090	U

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Mol	Chain	Res	Type
36	1	3091	A
36	1	3092	C
36	1	3093	C
36	1	3097	C
36	1	3120	C
36	1	3122	A
36	1	3129	A
36	1	3130	A
36	1	3131	U
36	1	3141	A
36	1	3142	A
36	1	3143	C
36	1	3145	C
36	1	3151	U
36	1	3153	U
36	1	3154	C
36	1	3155	U
36	1	3156	U
36	1	3157	U
36	1	3165	A
36	1	3167	A
36	1	3168	A
36	1	3170	A
36	1	3171	U
36	1	3173	G
36	1	3174	A
36	1	3176	G
36	1	3179	U
36	1	3181	C
36	1	3187	A
36	1	3196	U
36	1	3198	U
36	1	3207	U
36	1	3209	A
36	1	3210	A
36	1	3217	C
36	1	3218	A
36	1	3219	G
36	1	3223	A
36	1	3228	C
36	1	3229	G
36	1	3244	A

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Mol	Chain	Res	Type
36	1	3245	A
36	1	3246	G
36	1	3247	G
36	1	3259	U
36	1	3263	G
36	1	3269	U
36	1	3270	U
36	1	3271	G
36	1	3276	G
36	1	3281	U
36	1	3286	G
36	1	3287	U
36	1	3288	G
36	1	3289	G
36	1	3294	A
36	1	3295	A
36	1	3303	G
36	1	3304	U
36	1	3307	A
36	1	3309	G
36	1	3313	U
36	1	3316	A
36	1	3317	U
36	1	3318	G
36	1	3319	U
36	1	3332	U
36	1	3335	A
36	1	3341	U
36	1	3342	A
36	1	3345	G
36	1	3347	A
36	1	3351	U
36	1	3352	U
36	1	3354	U
36	1	3355	U
36	1	3356	G
36	1	3359	A
36	1	3360	C
36	1	3367	C
36	1	3368	U
36	1	3369	G
36	1	3378	C

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Mol	Chain	Res	Type
36	1	3382	U
36	1	3383	G
36	1	3389	U
36	1	3390	G
36	1	3396	U
37	3	7	G
37	3	22	A
37	3	41	G
37	3	65	G
37	3	68	C
37	3	74	C
37	3	76	A
37	3	95	A
37	3	102	A
37	3	112	G
37	3	121	U
38	4	34	U
38	4	35	C
38	4	48	A
38	4	53	A
38	4	59	A
38	4	60	U
38	4	62	C
38	4	63	G
38	4	75	G
38	4	80	A
38	4	81	U
38	4	83	C
38	4	85	G
38	4	86	U
38	4	87	G
38	4	90	U
38	4	95	G
38	4	104	A
38	4	105	A
38	4	106	C
38	4	107	G
38	4	111	A
38	4	112	U
38	4	113	U
38	4	125	U
38	4	126	A

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Mol	Chain	Res	Type
38	4	128	U
38	4	149	A
38	4	152	G
38	4	155	A
38	4	157	U
38	4	158	U
1	6	4	C
1	6	17	C
1	6	23	G
1	6	25	C
1	6	26	A
1	6	27	U
1	6	34	G
1	6	42	G
1	6	46	A
1	6	47	A
1	6	57	G
1	6	60	U
1	6	63	G
1	6	65	A
1	6	66	U
1	6	68	A
1	6	69	G
1	6	70	C
1	6	71	A
1	6	72	A
1	6	73	U
1	6	75	U
1	6	76	A
1	6	77	U
1	6	78	A
1	6	95	G
1	6	101	U
1	6	104	A
1	6	111	U
1	6	114	C
1	6	116	U
1	6	124	A
1	6	127	G
1	6	132	U
1	6	137	U
1	6	140	A

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Mol	Chain	Res	Type
1	6	141	U
1	6	144	U
1	6	145	A
1	6	146	U
1	6	153	G
1	6	155	U
1	6	158	U
1	6	159	U
1	6	166	C
1	6	178	U
1	6	179	A
1	6	181	A
1	6	185	U
1	6	188	A
1	6	191	C
1	6	192	U
1	6	193	U
1	6	194	U
1	6	195	G
1	6	196	G
1	6	199	G
1	6	200	A
1	6	217	A
1	6	218	A
1	6	219	A
1	6	220	A
1	6	223	U
1	6	224	C
1	6	227	U
1	6	228	G
1	6	230	C
1	6	232	U
1	6	233	C
1	6	234	G
1	6	240	U
1	6	241	U
1	6	245	U
1	6	250	C
1	6	260	U
1	6	261	U
1	6	265	A
1	6	266	A

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Mol	Chain	Res	Type
1	6	271	A
1	6	272	U
1	6	273	G
1	6	277	U
1	6	278	U
1	6	280	U
1	6	281	G
1	6	287	G
1	6	294	C
1	6	299	A
1	6	302	U
1	6	308	C
1	6	314	C
1	6	316	A
1	6	319	U
1	6	321	C
1	6	337	G
1	6	338	C
1	6	343	C
1	6	352	A
1	6	359	A
1	6	361	C
1	6	380	U
1	6	400	A
1	6	401	A
1	6	402	C
1	6	404	G
1	6	416	A
1	6	418	G
1	6	421	A
1	6	424	C
1	6	425	A
1	6	426	G
1	6	428	A
1	6	434	G
1	6	439	U
1	6	444	C
1	6	445	A
1	6	448	C
1	6	468	A
1	6	475	A
1	6	477	A

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Mol	Chain	Res	Type
1	6	484	C
1	6	486	G
1	6	487	G
1	6	488	G
1	6	489	C
1	6	490	C
1	6	492	A
1	6	493	U
1	6	494	U
1	6	496	G
1	6	500	C
1	6	501	U
1	6	505	A
1	6	506	A
1	6	508	U
1	6	509	G
1	6	511	A
1	6	516	G
1	6	519	C
1	6	528	U
1	6	534	A
1	6	536	C
1	6	539	G
1	6	541	A
1	6	542	A
1	6	543	C
1	6	544	A
1	6	548	G
1	6	554	C
1	6	555	A
1	6	556	A
1	6	557	G
1	6	558	U
1	6	559	C
1	6	564	G
1	6	565	C
1	6	566	C
1	6	570	A
1	6	574	G
1	6	579	A
1	6	580	A
1	6	582	U

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Mol	Chain	Res	Type
1	6	594	A
1	6	595	G
1	6	597	G
1	6	606	A
1	6	608	U
1	6	609	U
1	6	610	G
1	6	611	U
1	6	619	A
1	6	620	A
1	6	623	A
1	6	634	G
1	6	637	C
1	6	639	U
1	6	640	U
1	6	650	U
1	6	651	G
1	6	652	G
1	6	653	C
1	6	661	A
1	6	662	U
1	6	665	U
1	6	667	U
1	6	668	C
1	6	670	U
1	6	679	U
1	6	680	U
1	6	681	U
1	6	682	C
1	6	683	C
1	6	684	A
1	6	685	A
1	6	690	G
1	6	698	U
1	6	709	C
1	6	710	U
1	6	711	U
1	6	715	U
1	6	718	U
1	6	719	U
1	6	720	G
1	6	721	U

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Mol	Chain	Res	Type
1	6	722	G
1	6	723	G
1	6	730	G
1	6	742	U
1	6	751	G
1	6	754	A
1	6	755	A
1	6	756	A
1	6	765	G
1	6	770	A
1	6	774	A
1	6	775	G
1	6	780	A
1	6	781	U
1	6	782	U
1	6	783	G
1	6	787	G
1	6	793	A
1	6	794	U
1	6	801	G
1	6	803	A
1	6	809	A
1	6	811	A
1	6	812	A
1	6	814	A
1	6	815	G
1	6	821	U
1	6	823	G
1	6	824	G
1	6	825	U
1	6	826	U
1	6	829	A
1	6	830	U
1	6	831	U
1	6	832	U
1	6	834	G
1	6	835	U
1	6	861	U
1	6	862	A
1	6	863	A
1	6	865	A
1	6	876	G

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Mol	Chain	Res	Type
1	6	886	U
1	6	898	A
1	6	906	A
1	6	913	G
1	6	914	G
1	6	933	A
1	6	935	U
1	6	942	G
1	6	951	A
1	6	959	U
1	6	960	U
1	6	966	A
1	6	969	C
1	6	971	A
1	6	983	A
1	6	988	A
1	6	992	A
1	6	993	A
1	6	996	U
1	6	998	A
1	6	1000	C
1	6	1003	A
1	6	1004	U
1	6	1005	A
1	6	1016	C
1	6	1021	C
1	6	1026	A
1	6	1028	C
1	6	1029	U
1	6	1039	A
1	6	1040	G
1	6	1052	U
1	6	1053	G
1	6	1057	U
1	6	1058	U
1	6	1059	U
1	6	1060	U
1	6	1072	C
1	6	1075	C
1	6	1081	A
1	6	1082	C
1	6	1092	A

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Mol	Chain	Res	Type
1	6	1096	C
1	6	1097	U
1	6	1098	U
1	6	1100	G
1	6	1101	G
1	6	1109	G
1	6	1111	G
1	6	1137	A
1	6	1138	A
1	6	1151	A
1	6	1155	G
1	6	1158	C
1	6	1159	C
1	6	1160	A
1	6	1161	C
1	6	1167	G
1	6	1183	A
1	6	1185	U
1	6	1191	U
1	6	1194	A
1	6	1196	A
1	6	1199	G
1	6	1200	G
1	6	1202	A
1	6	1207	C
1	6	1217	A
1	6	1218	G
1	6	1225	U
1	6	1226	A
1	6	1228	G
1	6	1229	G
1	6	1230	A
1	6	1239	U
1	6	1241	G
1	6	1242	A
1	6	1243	G
1	6	1245	G
1	6	1246	C
1	6	1252	C
1	6	1256	A
1	6	1257	U
1	6	1258	U

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Mol	Chain	Res	Type
1	6	1284	C
1	6	1286	U
1	6	1288	G
1	6	1291	G
1	6	1314	U
1	6	1315	U
1	6	1316	G
1	6	1318	G
1	6	1321	A
1	6	1335	U
1	6	1338	C
1	6	1344	A
1	6	1345	A
1	6	1346	A
1	6	1353	U
1	6	1361	U
1	6	1363	U
1	6	1364	G
1	6	1367	G
1	6	1370	U
1	6	1371	A
1	6	1383	G
1	6	1390	U
1	6	1398	U
1	6	1399	C
1	6	1400	A
1	6	1401	A
1	6	1402	G
1	6	1413	U
1	6	1414	U
1	6	1415	U
1	6	1418	G
1	6	1426	C
1	6	1427	A
1	6	1428	G
1	6	1445	G
1	6	1446	A
1	6	1447	C
1	6	1448	G
1	6	1458	G
1	6	1459	C
1	6	1460	A

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Mol	Chain	Res	Type
1	6	1461	C
1	6	1466	G
1	6	1471	A
1	6	1473	U
1	6	1481	C
1	6	1482	C
1	6	1486	G
1	6	1490	C
1	6	1491	U
1	6	1492	A
1	6	1493	A
1	6	1494	C
1	6	1504	G
1	6	1506	G
1	6	1514	U
1	6	1516	A
1	6	1519	U
1	6	1521	G
1	6	1522	U
1	6	1523	G
1	6	1524	A
1	6	1535	U
1	6	1536	G
1	6	1537	C
1	6	1538	U
1	6	1540	G
1	6	1554	U
1	6	1557	U
1	6	1559	A
1	6	1573	A
1	6	1574	G
1	6	1575	G
1	6	1579	U
1	6	1582	U
1	6	1584	G
1	6	1600	A
1	6	1601	G
1	6	1621	U
1	6	1631	A
1	6	1634	C
1	6	1636	C
1	6	1637	C

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Mol	Chain	Res	Type
1	6	1638	G
1	6	1657	U
1	6	1658	G
1	6	1668	G
1	6	1683	C
1	6	1696	G
1	6	1698	G
1	6	1699	G
1	6	1700	C
1	6	1701	A
1	6	1702	A
1	6	1712	A
1	6	1713	G
1	6	1716	C
1	6	1717	G
1	6	1731	A
1	6	1736	G
1	6	1742	U
1	6	1760	G
1	6	1762	A
1	6	1766	A
1	6	1767	G
1	6	1769	U
1	6	1780	G
1	6	1782	A
1	6	1789	G
1	6	1792	G
1	6	1793	G
1	6	1794	A
1	6	1795	U
1	6	1796	C
1	6	1799	U
1	6	1800	A
36	5	6	A
36	5	15	C
36	5	16	A
36	5	26	A
36	5	30	G
36	5	40	A
36	5	43	A
36	5	49	A
36	5	57	A

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Mol	Chain	Res	Type
36	5	59	G
36	5	60	A
36	5	65	A
36	5	66	A
36	5	71	A
36	5	73	C
36	5	76	G
36	5	92	G
36	5	96	G
36	5	109	A
36	5	110	G
36	5	111	C
36	5	113	C
36	5	118	U
36	5	121	A
36	5	122	A
36	5	131	C
36	5	134	U
36	5	135	C
36	5	136	G
36	5	156	G
36	5	157	A
36	5	165	A
36	5	168	U
36	5	170	G
36	5	171	G
36	5	172	G
36	5	173	G
36	5	174	C
36	5	180	C
36	5	182	U
36	5	187	A
36	5	190	U
36	5	191	U
36	5	196	G
36	5	197	G
36	5	200	C
36	5	201	A
36	5	210	U
36	5	213	A
36	5	218	G
36	5	219	A

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Mol	Chain	Res	Type
36	5	220	G
36	5	221	A
36	5	231	G
36	5	235	A
36	5	239	G
36	5	240	U
36	5	241	G
36	5	245	U
36	5	247	C
36	5	248	U
36	5	249	U
36	5	250	U
36	5	252	U
36	5	253	A
36	5	254	A
36	5	269	G
36	5	270	U
36	5	284	A
36	5	286	U
36	5	295	A
36	5	298	U
36	5	305	U
36	5	311	C
36	5	315	C
36	5	323	A
36	5	326	U
36	5	327	A
36	5	329	U
36	5	334	A
36	5	339	C
36	5	349	A
36	5	350	C
36	5	351	A
36	5	360	G
36	5	376	G
36	5	379	C
36	5	390	G
36	5	395	A
36	5	398	A
36	5	399	A
36	5	401	U
36	5	402	A

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Mol	Chain	Res	Type
36	5	403	C
36	5	404	G
36	5	418	A
36	5	421	G
36	5	422	A
36	5	436	A
36	5	437	G
36	5	439	C
36	5	440	A
36	5	441	U
36	5	442	G
36	5	492	U
36	5	496	C
36	5	521	A
36	5	542	G
36	5	546	C
36	5	547	G
36	5	548	G
36	5	550	A
36	5	551	A
36	5	554	A
36	5	557	A
36	5	559	A
36	5	578	A
36	5	579	G
36	5	581	U
36	5	592	A
36	5	600	G
36	5	609	G
36	5	611	A
36	5	619	A
36	5	621	A
36	5	636	C
36	5	647	A
36	5	649	A
36	5	651	G
36	5	660	A
36	5	662	U
36	5	677	A
36	5	681	U
36	5	691	A
36	5	692	A

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Mol	Chain	Res	Type
36	5	700	C
36	5	705	A
36	5	712	G
36	5	715	A
36	5	716	A
36	5	725	G
36	5	736	A
36	5	766	U
36	5	767	U
36	5	776	U
36	5	777	U
36	5	781	G
36	5	785	G
36	5	786	A
36	5	806	A
36	5	811	U
36	5	817	A
36	5	826	G
36	5	830	A
36	5	847	A
36	5	861	C
36	5	874	U
36	5	879	U
36	5	881	C
36	5	890	C
36	5	907	G
36	5	908	G
36	5	914	A
36	5	916	G
36	5	917	A
36	5	921	A
36	5	923	C
36	5	924	G
36	5	925	A
36	5	937	G
36	5	944	C
36	5	959	C
36	5	960	U
36	5	962	A
36	5	963	G
36	5	964	G
36	5	974	G

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Mol	Chain	Res	Type
36	5	979	U
36	5	994	G
36	5	1000	C
36	5	1001	G
36	5	1002	A
36	5	1010	G
36	5	1015	U
36	5	1016	C
36	5	1017	C
36	5	1018	G
36	5	1021	G
36	5	1024	G
36	5	1025	A
36	5	1026	A
36	5	1028	U
36	5	1029	G
36	5	1035	G
36	5	1047	A
36	5	1049	C
36	5	1063	G
36	5	1064	A
36	5	1065	A
36	5	1072	G
36	5	1081	U
36	5	1082	U
36	5	1085	A
36	5	1087	G
36	5	1093	A
36	5	1095	U
36	5	1097	G
36	5	1098	A
36	5	1103	A
36	5	1104	G
36	5	1117	G
36	5	1131	G
36	5	1143	A
36	5	1144	U
36	5	1152	G
36	5	1153	A
36	5	1159	A
36	5	1160	C
36	5	1169	A

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Mol	Chain	Res	Type
36	5	1175	C
36	5	1177	G
36	5	1178	G
36	5	1179	A
36	5	1180	A
36	5	1181	U
36	5	1191	U
36	5	1196	C
36	5	1201	C
36	5	1202	A
36	5	1206	G
36	5	1209	G
36	5	1222	G
36	5	1223	A
36	5	1235	U
36	5	1236	G
36	5	1239	C
36	5	1241	U
36	5	1242	G
36	5	1243	G
36	5	1245	A
36	5	1246	G
36	5	1252	A
36	5	1253	U
36	5	1258	U
36	5	1262	G
36	5	1263	A
36	5	1264	G
36	5	1265	U
36	5	1266	G
36	5	1277	C
36	5	1307	G
36	5	1308	A
36	5	1309	U
36	5	1313	G
36	5	1316	C
36	5	1330	A
36	5	1331	U
36	5	1332	A
36	5	1348	U
36	5	1349	G
36	5	1351	U

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Mol	Chain	Res	Type
36	5	1353	U
36	5	1354	G
36	5	1355	A
36	5	1356	U
36	5	1357	G
36	5	1386	A
36	5	1387	G
36	5	1391	C
36	5	1398	U
36	5	1399	A
36	5	1400	G
36	5	1408	G
36	5	1417	G
36	5	1418	A
36	5	1419	A
36	5	1421	G
36	5	1428	A
36	5	1431	G
36	5	1433	A
36	5	1434	G
36	5	1437	C
36	5	1438	U
36	5	1446	A
36	5	1450	G
36	5	1471	U
36	5	1481	A
36	5	1482	A
36	5	1490	A
36	5	1503	A
36	5	1508	C
36	5	1523	U
36	5	1527	C
36	5	1536	G
36	5	1554	U
36	5	1555	U
36	5	1556	C
36	5	1557	A
36	5	1560	G
36	5	1561	G
36	5	1562	C
36	5	1567	U
36	5	1569	U

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Mol	Chain	Res	Type
36	5	1570	U
36	5	1571	A
36	5	1574	C
36	5	1575	A
36	5	1576	G
36	5	1577	G
36	5	1578	C
36	5	1579	C
36	5	1581	C
36	5	1582	C
36	5	1583	A
36	5	1587	A
36	5	1589	A
36	5	1593	A
36	5	1605	A
36	5	1620	U
36	5	1621	A
36	5	1629	U
36	5	1632	A
36	5	1639	C
36	5	1643	A
36	5	1644	C
36	5	1645	U
36	5	1657	C
36	5	1677	G
36	5	1683	A
36	5	1687	U
36	5	1694	U
36	5	1702	U
36	5	1714	A
36	5	1716	U
36	5	1717	U
36	5	1724	U
36	5	1725	C
36	5	1730	G
36	5	1750	A
36	5	1751	G
36	5	1760	A
36	5	1762	C
36	5	1764	U
36	5	1765	U
36	5	1766	G

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Mol	Chain	Res	Type
36	5	1769	G
36	5	1770	G
36	5	1775	G
36	5	1780	G
36	5	1797	A
36	5	1812	G
36	5	1814	A
36	5	1815	U
36	5	1816	A
36	5	1817	G
36	5	1818	U
36	5	1821	U
36	5	1829	G
36	5	1839	A
36	5	1841	A
36	5	1842	A
36	5	1846	C
36	5	1847	A
36	5	1849	C
36	5	1850	A
36	5	1878	G
36	5	1879	A
36	5	1880	U
36	5	1895	A
36	5	1901	A
36	5	1906	G
36	5	1935	G
36	5	2101	C
36	5	2102	U
36	5	2111	G
36	5	2112	U
36	5	2113	A
36	5	2115	G
36	5	2121	G
36	5	2122	G
36	5	2131	A
36	5	2144	A
36	5	2149	A
36	5	2157	G
36	5	2158	A
36	5	2168	A
36	5	2169	G

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Mol	Chain	Res	Type
36	5	2187	G
36	5	2188	A
36	5	2205	U
36	5	2206	G
36	5	2210	G
36	5	2215	A
36	5	2223	A
36	5	2225	U
36	5	2231	C
36	5	2244	A
36	5	2246	G
36	5	2250	G
36	5	2253	G
36	5	2255	A
36	5	2256	A
36	5	2258	U
36	5	2267	C
36	5	2269	U
36	5	2270	A
36	5	2272	G
36	5	2273	G
36	5	2276	G
36	5	2279	A
36	5	2281	A
36	5	2288	G
36	5	2298	U
36	5	2303	A
36	5	2307	G
36	5	2308	C
36	5	2310	U
36	5	2313	A
36	5	2315	G
36	5	2318	U
36	5	2335	G
36	5	2336	U
36	5	2364	G
36	5	2367	A
36	5	2372	A
36	5	2373	A
36	5	2374	C
36	5	2375	G
36	5	2385	G

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Mol	Chain	Res	Type
36	5	2393	G
36	5	2397	A
36	5	2401	A
36	5	2403	G
36	5	2404	A
36	5	2411	U
36	5	2418	G
36	5	2419	A
36	5	2435	G
36	5	2439	A
36	5	2446	U
36	5	2447	A
36	5	2452	G
36	5	2494	A
36	5	2497	U
36	5	2501	U
36	5	2502	A
36	5	2503	G
36	5	2504	U
36	5	2505	U
36	5	2514	U
36	5	2515	A
36	5	2518	C
36	5	2523	A
36	5	2529	A
36	5	2533	G
36	5	2536	A
36	5	2537	U
36	5	2538	U
36	5	2539	C
36	5	2540	A
36	5	2543	U
36	5	2549	G
36	5	2552	C
36	5	2555	G
36	5	2560	C
36	5	2564	G
36	5	2565	U
36	5	2567	C
36	5	2568	C
36	5	2569	A
36	5	2570	U

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Mol	Chain	Res	Type
36	5	2571	U
36	5	2572	C
36	5	2573	G
36	5	2574	G
36	5	2584	G
36	5	2585	G
36	5	2589	G
36	5	2593	A
36	5	2606	G
36	5	2607	G
36	5	2614	G
36	5	2617	U
36	5	2626	A
36	5	2637	A
36	5	2652	U
36	5	2656	A
36	5	2674	A
36	5	2676	A
36	5	2677	G
36	5	2678	A
36	5	2689	A
36	5	2691	A
36	5	2694	A
36	5	2696	A
36	5	2704	A
36	5	2705	A
36	5	2714	G
36	5	2717	U
36	5	2727	A
36	5	2728	G
36	5	2729	U
36	5	2737	C
36	5	2748	A
36	5	2752	U
36	5	2753	G
36	5	2755	C
36	5	2762	A
36	5	2772	C
36	5	2773	C
36	5	2777	G
36	5	2778	G
36	5	2780	A

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Mol	Chain	Res	Type
36	5	2796	G
36	5	2797	C
36	5	2799	A
36	5	2800	G
36	5	2801	A
36	5	2810	C
36	5	2814	G
36	5	2816	G
36	5	2817	A
36	5	2818	U
36	5	2829	U
36	5	2838	A
36	5	2843	U
36	5	2845	A
36	5	2847	A
36	5	2853	A
36	5	2855	U
36	5	2861	U
36	5	2862	U
36	5	2863	G
36	5	2871	G
36	5	2886	U
36	5	2887	A
36	5	2898	G
36	5	2899	C
36	5	2904	U
36	5	2923	U
36	5	2935	U
36	5	2936	A
36	5	2942	C
36	5	2945	G
36	5	2947	G
36	5	2954	U
36	5	2957	G
36	5	2965	U
36	5	2971	A
36	5	2977	G
36	5	2983	C
36	5	2990	G
36	5	2996	U
36	5	2997	G
36	5	3012	A

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Mol	Chain	Res	Type
36	5	3018	C
36	5	3030	G
36	5	3056	U
36	5	3059	G
36	5	3078	U
36	5	3079	U
36	5	3086	A
36	5	3087	A
36	5	3091	A
36	5	3092	C
36	5	3102	G
36	5	3119	U
36	5	3122	A
36	5	3130	A
36	5	3131	U
36	5	3142	A
36	5	3143	C
36	5	3150	A
36	5	3152	U
36	5	3153	U
36	5	3155	U
36	5	3156	U
36	5	3157	U
36	5	3158	G
36	5	3164	C
36	5	3165	A
36	5	3168	A
36	5	3171	U
36	5	3172	A
36	5	3173	G
36	5	3174	A
36	5	3176	G
36	5	3177	G
36	5	3179	U
36	5	3180	A
36	5	3181	C
36	5	3187	A
36	5	3195	U
36	5	3196	U
36	5	3197	G
36	5	3207	U
36	5	3217	C

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Mol	Chain	Res	Type
36	5	3218	A
36	5	3219	G
36	5	3224	G
36	5	3229	G
36	5	3234	A
36	5	3238	G
36	5	3239	G
36	5	3243	A
36	5	3244	A
36	5	3245	A
36	5	3246	G
36	5	3247	G
36	5	3258	U
36	5	3259	U
36	5	3275	U
36	5	3276	G
36	5	3277	U
36	5	3281	U
36	5	3282	U
36	5	3284	G
36	5	3285	C
36	5	3286	G
36	5	3288	G
36	5	3289	G
36	5	3290	G
36	5	3294	A
36	5	3304	U
36	5	3305	A
36	5	3313	U
36	5	3316	A
36	5	3317	U
36	5	3318	G
36	5	3320	A
36	5	3342	A
36	5	3345	G
36	5	3348	G
36	5	3351	U
36	5	3352	U
36	5	3353	G
36	5	3354	U
36	5	3358	U
36	5	3361	G

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Mol	Chain	Res	Type
36	5	3369	G
36	5	3377	G
36	5	3378	C
36	5	3381	U
36	5	3382	U
36	5	3389	U
36	5	3390	G
36	5	3393	U
36	5	3394	U
36	5	3396	U
37	7	7	G
37	7	22	A
37	7	41	G
37	7	47	C
37	7	54	U
37	7	55	A
37	7	60	G
37	7	65	G
37	7	73	C
37	7	76	A
37	7	99	G
37	7	102	A
37	7	112	G
38	8	34	U
38	8	35	C
38	8	48	A
38	8	53	A
38	8	57	C
38	8	58	G
38	8	59	A
38	8	60	U
38	8	62	C
38	8	63	G
38	8	75	G
38	8	79	A
38	8	80	A
38	8	81	U
38	8	82	U
38	8	83	C
38	8	84	C
38	8	86	U
38	8	87	G

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Mol	Chain	Res	Type
38	8	95	G
38	8	96	A
38	8	97	A
38	8	99	C
38	8	104	A
38	8	106	C
38	8	111	A
38	8	113	U
38	8	125	U
38	8	126	A
38	8	127	U
38	8	129	C
38	8	152	G
38	8	157	U
38	8	158	U

All (177) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	2	25	C
1	2	114	C
1	2	130	C
1	2	131	C
1	2	158	U
1	2	218	A
1	2	240	U
1	2	278	U
1	2	280	U
1	2	417	A
1	2	497	G
1	2	499	U
1	2	501	U
1	2	503	G
1	2	512	A
1	2	558	U
1	2	622	A
1	2	685	A
1	2	704	C
1	2	720	G
1	2	721	U
1	2	755	A
1	2	829	A

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Mol	Chain	Res	Type
1	2	1081	A
1	2	1150	G
1	2	1196	A
1	2	1207	C
1	2	1226	A
1	2	1244	A
1	2	1250	U
1	2	1344	A
1	2	1370	U
1	2	1481	C
1	2	1489	U
1	2	1568	C
1	2	1573	A
1	2	1615	C
1	2	1657	U
1	2	1698	G
1	2	1761	U
36	1	65	A
36	1	210	U
36	1	239	G
36	1	282	G
36	1	547	G
36	1	659	G
36	1	715	A
36	1	763	G
36	1	873	C
36	1	916	G
36	1	981	U
36	1	993	G
36	1	1064	A
36	1	1094	U
36	1	1097	G
36	1	1103	A
36	1	1181	U
36	1	1196	C
36	1	1307	G
36	1	1329	U
36	1	1331	U
36	1	1352	A
36	1	1355	A
36	1	1484	U
36	1	1507	G

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Mol	Chain	Res	Type
36	1	1562	C
36	1	1582	C
36	1	1716	U
36	1	1820	U
36	1	2101	C
36	1	2112	U
36	1	2209	U
36	1	2249	G
36	1	2297	U
36	1	2372	A
36	1	2400	G
36	1	2418	G
36	1	2537	U
36	1	2541	U
36	1	2554	A
36	1	2585	G
36	1	2593	A
36	1	2817	A
36	1	2818	U
36	1	2971	A
36	1	3078	U
36	1	3121	U
36	1	3228	C
36	1	3269	U
36	1	3350	C
36	1	3351	U
36	1	3353	G
36	1	3377	G
38	4	80	A
38	4	85	G
38	4	125	U
1	6	25	C
1	6	114	C
1	6	158	U
1	6	187	G
1	6	192	U
1	6	217	A
1	6	259	U
1	6	400	A
1	6	417	A
1	6	454	U
1	6	542	A

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Mol	Chain	Res	Type
1	6	555	A
1	6	558	U
1	6	697	C
1	6	717	C
1	6	720	G
1	6	755	A
1	6	828	U
1	6	1058	U
1	6	1097	U
1	6	1244	A
1	6	1255	G
1	6	1344	A
1	6	1481	C
1	6	1491	U
1	6	1535	U
1	6	1573	A
1	6	1600	A
1	6	1620	C
36	5	210	U
36	5	238	A
36	5	438	A
36	5	594	U
36	5	873	C
36	5	916	G
36	5	978	G
36	5	993	G
36	5	1027	A
36	5	1064	A
36	5	1081	U
36	5	1152	G
36	5	1190	A
36	5	1222	G
36	5	1238	C
36	5	1241	U
36	5	1307	G
36	5	1329	U
36	5	1331	U
36	5	1352	A
36	5	1355	A
36	5	1481	A
36	5	1560	G
36	5	1580	A

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Mol	Chain	Res	Type
36	5	1716	U
36	5	1816	A
36	5	1841	A
36	5	1842	A
36	5	2101	C
36	5	2112	U
36	5	2209	U
36	5	2249	G
36	5	2255	A
36	5	2257	C
36	5	2372	A
36	5	2400	G
36	5	2418	G
36	5	2500	A
36	5	2513	U
36	5	2539	C
36	5	2772	C
36	5	2818	U
36	5	2872	A
36	5	3078	U
36	5	3121	U
36	5	3195	U
36	5	3207	U
36	5	3218	A
36	5	3275	U
36	5	3289	G
36	5	3317	U
36	5	3357	U

## 5.4 Non-standard residues in protein, DNA, RNA chains

6 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
84	CH	B	74	84	14,21,22	0.82	0	18,30,33	0.49	0
84	CH	B	75	87,84	14,21,22	0.66	0	18,30,33	0.82	0
84	8AN	B	76	87,84	18,24,25	1.11	1 (5%)	10,35,38	2.14	2 (20%)
84	CH	C	74	84	15,18,22	0.79	1 (6%)	18,26,33	0.75	0
84	CH	C	75	87,84	14,21,22	0.74	0	18,30,33	0.54	0
84	8AN	C	76	87,84	18,24,25	1.14	1 (5%)	10,35,38	2.12	1 (10%)

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
84	CH	B	74	84	-	0/3/25/26	0/2/2/2
84	CH	B	75	87,84	-	0/3/25/26	0/2/2/2
84	8AN	B	76	87,84	-	0/3/25/26	0/3/3/3
84	CH	C	74	84	-	0/2/22/26	0/2/2/2
84	CH	C	75	87,84	-	0/3/25/26	0/2/2/2
84	8AN	C	76	87,84	-	0/3/25/26	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
84	C	74	CH	C6-C5	-2.03	1.33	1.38
84	C	76	8AN	C5-C4	2.93	1.47	1.40
84	B	76	8AN	C5-C4	3.25	1.47	1.40

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
84	C	76	8AN	N3-C2-N1	-5.83	124.30	128.87
84	B	76	8AN	N3-C2-N1	-5.71	124.39	128.87
84	B	76	8AN	N6-C6-N1	2.18	122.18	118.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
84	B	74	CH	1	0
84	B	75	CH	1	0
84	B	76	8AN	1	0
84	C	74	CH	3	0
84	C	75	CH	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 33 ligands modelled in this entry, 28 are monoatomic - leaving 5 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$
86	SPS	1	3401	-	19,23,23	3.50	10 (52%)	16,30,30	2.99	7 (43%)
88	OHX	1	3405	-	0,6,6	0.00	-	0,15,15	0.00	-
88	OHX	1	3406	-	0,6,6	0.00	-	0,15,15	0.00	-
88	OHX	5	3406	-	0,6,6	0.00	-	0,15,15	0.00	-
86	SPS	B	3401	87	19,23,23	3.49	11 (57%)	16,30,30	3.14	8 (50%)

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
86	SPS	1	3401	-	-	0/15/18/18	0/1/1/1
88	OHX	1	3405	-	-	0/0/0/0	0/0/0/0
88	OHX	1	3406	-	-	0/0/0/0	0/0/0/0
88	OHX	5	3406	-	-	0/0/0/0	0/0/0/0
86	SPS	B	3401	87	-	0/15/18/18	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
86	1	3401	SPS	C9-C10	-8.75	1.31	1.48
86	B	3401	SPS	C9-C10	-8.45	1.31	1.48
86	B	3401	SPS	O13-C13	-5.31	1.19	1.42
86	1	3401	SPS	O13-C13	-5.20	1.19	1.42
86	B	3401	SPS	O10-C10	-2.99	1.18	1.24
86	1	3401	SPS	O10-C10	-2.76	1.19	1.24
86	B	3401	SPS	C1-C6	-2.60	1.38	1.44
86	B	3401	SPS	C3-N4	-2.33	1.33	1.38
86	B	3401	SPS	C3-N2	-2.25	1.33	1.38
86	1	3401	SPS	C1-C6	-2.23	1.39	1.44
86	1	3401	SPS	O1-C1	-2.13	1.19	1.24
86	B	3401	SPS	O1-C1	-2.04	1.19	1.24
86	1	3401	SPS	C3-N4	-2.03	1.33	1.38
86	B	3401	SPS	C6-C8	2.92	1.53	1.47
86	1	3401	SPS	C6-C8	3.24	1.54	1.47
86	B	3401	SPS	C10-N11	3.30	1.44	1.34
86	1	3401	SPS	C10-N11	3.71	1.45	1.34
86	B	3401	SPS	O15-S15	4.40	1.64	1.50
86	1	3401	SPS	O15-S15	4.66	1.65	1.50
86	1	3401	SPS	C9-C8	7.05	1.52	1.32
86	B	3401	SPS	C9-C8	7.33	1.53	1.32

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
86	B	3401	SPS	N2-C3-N4	-5.15	119.01	127.69
86	1	3401	SPS	C7-C5-C6	-4.46	118.29	123.63
86	1	3401	SPS	N2-C3-N4	-4.33	120.39	127.69
86	B	3401	SPS	C7-C5-C6	-2.82	120.25	123.63
86	1	3401	SPS	C6-C8-C9	-2.77	118.07	127.08
86	B	3401	SPS	C12-N11-C10	-2.75	118.86	122.57
86	B	3401	SPS	C6-C8-C9	-2.61	118.59	127.08
86	B	3401	SPS	O13-C13-C12	2.51	119.20	112.24
86	B	3401	SPS	C9-C10-N11	2.56	119.86	114.14
86	1	3401	SPS	C18-S17-C16	2.57	108.02	100.14
86	1	3401	SPS	O13-C13-C12	2.83	120.08	112.24
86	1	3401	SPS	C7-C5-N4	2.93	120.77	116.32
86	B	3401	SPS	C18-S17-C16	3.59	111.12	100.14
86	1	3401	SPS	C1-N2-C3	7.67	121.56	115.16
86	B	3401	SPS	C1-N2-C3	8.39	122.15	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
86	1	3401	SPS	2	0
86	B	3401	SPS	3	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
80	m2	2
1	2	2
35	SM	1
35	sM	1
12	c0	1
36	5	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	sM	139:UNK	C	155:UNK	N	38.68
1	SM	141:ALA	C	151:UNK	N	26.53
1	c0	84:UNK	C	87:UNK	N	8.84
1	2	1716:C	O3'	1717:G	P	5.83
1	5	2437:G	O3'	2438:A	P	4.00
1	m2	23:UNK	C	28:UNK	N	3.70
1	m2	52:UNK	C	54:UNK	N	3.66
1	2	1685:G	O3'	1686:C	P	3.03

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	2	1781/1800 (98%)	0.48	91 (5%) 32 25	60, 98, 170, 210	0
1	6	1795/1800 (99%)	0.27	58 (3%) 51 44	44, 86, 157, 204	0
2	S0	206/251 (82%)	1.22	49 (23%) 1 1	100, 112, 123, 128	0
2	s0	206/251 (82%)	1.16	50 (24%) 1 1	80, 97, 111, 117	0
3	S1	214/254 (84%)	1.60	77 (35%) 0 1	110, 145, 169, 175	0
3	s1	216/254 (85%)	1.03	44 (20%) 1 1	80, 96, 114, 123	0
4	S2	217/253 (85%)	0.73	30 (13%) 4 3	83, 96, 109, 113	0
4	s2	217/253 (85%)	0.45	9 (4%) 41 34	68, 79, 93, 108	0
5	S3	223/239 (93%)	0.67	26 (11%) 6 5	91, 100, 123, 128	0
5	s3	223/239 (93%)	1.46	72 (32%) 1 1	86, 112, 131, 137	0
6	S4	260/260 (100%)	1.31	67 (25%) 1 1	77, 99, 107, 124	0
6	s4	260/260 (100%)	1.20	60 (23%) 1 1	60, 82, 92, 119	0
7	S5	206/224 (91%)	1.21	41 (19%) 1 1	104, 124, 131, 134	0
7	s5	206/224 (91%)	0.62	21 (10%) 9 7	87, 108, 123, 130	0
8	S6	226/236 (95%)	0.85	41 (18%) 2 1	77, 105, 126, 132	0
8	s6	218/236 (92%)	0.49	21 (9%) 10 9	60, 89, 106, 114	0
9	S7	184/189 (97%)	0.90	30 (16%) 2 2	102, 121, 136, 142	0
9	s7	186/189 (98%)	0.67	23 (12%) 5 4	77, 104, 130, 134	0
10	S8	188/200 (94%)	0.78	21 (11%) 7 5	70, 86, 119, 129	0
10	s8	188/200 (94%)	1.01	28 (14%) 3 2	56, 73, 113, 130	0
11	S9	185/196 (94%)	1.69	71 (38%) 0 1	89, 106, 130, 141	0
11	s9	185/196 (94%)	1.21	44 (23%) 1 1	71, 88, 115, 126	0
12	C0	83/96 (86%)	2.34	51 (61%) 0 0	94, 112, 122, 126	0
12	c0	83/96 (86%)	2.04	40 (48%) 0 0	103, 131, 145, 149	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
13	C1	144/155 (92%)	1.04	21 (14%) 3 3	73, 83, 101, 113	0
13	c1	144/155 (92%)	0.84	9 (6%) 23 19	59, 70, 91, 113	0
14	C2	108/142 (76%)	0.52	14 (12%) 5 3	137, 145, 150, 152	0
14	c2	108/142 (76%)	0.82	16 (14%) 3 3	167, 183, 196, 199	0
15	C3	150/150 (100%)	0.59	16 (10%) 8 6	82, 95, 111, 115	0
15	c3	150/150 (100%)	0.14	5 (3%) 50 43	67, 79, 95, 100	0
16	C4	127/136 (93%)	1.42	35 (27%) 1 1	84, 142, 156, 161	0
16	c4	128/136 (94%)	1.53	43 (33%) 0 1	67, 101, 113, 116	0
17	C5	122/141 (86%)	0.94	21 (17%) 2 2	88, 99, 111, 117	0
17	c5	129/141 (91%)	0.67	16 (12%) 5 4	86, 103, 113, 119	0
18	C6	141/142 (99%)	1.70	48 (34%) 0 1	91, 112, 117, 120	0
18	c6	142/142 (100%)	0.82	24 (16%) 2 2	83, 102, 115, 126	0
19	C7	120/136 (88%)	1.44	33 (27%) 1 1	97, 109, 125, 129	0
19	c7	117/136 (86%)	0.38	10 (8%) 13 10	88, 101, 115, 119	0
20	C8	145/145 (100%)	0.95	24 (16%) 2 2	85, 110, 131, 138	0
20	c8	145/145 (100%)	0.80	23 (15%) 3 2	86, 101, 116, 126	0
21	C9	143/143 (100%)	1.40	37 (25%) 1 1	96, 112, 122, 126	0
21	c9	143/143 (100%)	1.63	52 (36%) 0 1	85, 97, 112, 118	0
22	D0	107/120 (89%)	0.98	20 (18%) 2 1	86, 112, 123, 125	0
22	d0	110/120 (91%)	2.37	58 (52%) 0 0	84, 114, 138, 147	0
23	D1	87/87 (100%)	0.91	16 (18%) 2 1	96, 104, 119, 127	0
23	d1	87/87 (100%)	0.70	9 (10%) 9 7	76, 88, 104, 112	0
24	D2	129/129 (100%)	0.64	11 (8%) 13 10	83, 95, 101, 109	0
24	d2	129/129 (100%)	0.46	9 (6%) 19 16	64, 74, 81, 87	0
25	D3	144/144 (100%)	0.81	18 (12%) 5 4	69, 77, 86, 96	0
25	d3	144/144 (100%)	0.60	7 (4%) 33 27	56, 61, 71, 79	0
26	D4	134/134 (100%)	1.30	33 (24%) 1 1	86, 107, 118, 129	0
26	d4	134/134 (100%)	0.36	10 (7%) 17 14	68, 90, 105, 110	0
27	D5	70/107 (65%)	1.51	24 (34%) 0 1	119, 130, 139, 140	0
27	d5	69/107 (64%)	1.51	19 (27%) 1 1	99, 113, 123, 125	0
28	D6	97/97 (100%)	2.01	44 (45%) 0 0	88, 106, 154, 159	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å <sup>2</sup> )	Q<0.9
28	d6	97/97 (100%)	1.37	26 (26%)	1	1	70, 87, 116, 123	0
29	D7	81/81 (100%)	1.23	17 (20%)	1	1	98, 110, 135, 141	0
29	d7	81/81 (100%)	0.94	12 (14%)	3	3	80, 93, 121, 128	0
30	D8	63/66 (95%)	1.21	15 (23%)	1	1	111, 127, 139, 143	0
30	d8	63/66 (95%)	1.06	10 (15%)	3	2	104, 116, 127, 129	0
31	D9	53/55 (96%)	1.26	12 (22%)	1	1	84, 91, 107, 112	0
31	d9	53/55 (96%)	2.26	26 (49%)	0	0	82, 92, 128, 139	0
32	E0	60/62 (96%)	1.84	21 (35%)	0	1	78, 107, 129, 134	0
32	e0	62/62 (100%)	0.83	11 (17%)	2	1	63, 88, 109, 122	0
33	E1	71/76 (93%)	0.66	9 (12%)	5	4	107, 133, 145, 148	0
33	e1	76/76 (100%)	1.09	20 (26%)	1	1	112, 158, 170, 175	0
34	SR	318/318 (100%)	1.49	102 (32%)	1	1	109, 120, 132, 146	0
34	sR	318/318 (100%)	1.56	105 (33%)	0	1	110, 126, 140, 150	0
35	SM	121/273 (44%)	0.86	22 (18%)	2	1	59, 94, 111, 116	0
35	sM	63/273 (23%)	1.60	21 (33%)	0	1	55, 105, 111, 112	0
36	1	3149/3396 (92%)	0.18	48 (1%)	76	71	35, 59, 123, 212	0
36	5	3169/3396 (93%)	0.21	48 (1%)	76	71	35, 54, 124, 188	0
37	3	121/121 (100%)	-0.00	0	100	100	43, 76, 90, 96	0
37	7	121/121 (100%)	-0.09	0	100	100	39, 59, 70, 75	0
38	4	158/158 (100%)	0.12	2 (1%)	79	74	43, 61, 91, 130	0
38	8	158/158 (100%)	0.11	1 (0%)	90	88	43, 63, 96, 120	0
39	L2	252/253 (99%)	0.36	9 (3%)	46	39	42, 61, 77, 82	0
39	l2	252/253 (99%)	0.41	11 (4%)	38	31	43, 59, 74, 83	0
40	L3	386/386 (100%)	0.45	19 (4%)	33	27	42, 63, 77, 85	0
40	l3	386/386 (100%)	0.33	8 (2%)	67	60	34, 49, 62, 77	0
41	L4	361/361 (100%)	-0.03	0	100	100	40, 54, 70, 71	0
41	l4	361/361 (100%)	0.14	2 (0%)	90	88	40, 57, 73, 88	0
42	L5	296/296 (100%)	1.15	68 (22%)	1	1	58, 81, 97, 111	0
42	l5	294/296 (99%)	0.61	21 (7%)	19	15	49, 60, 84, 93	0
43	L6	156/175 (89%)	0.59	12 (7%)	16	13	49, 57, 72, 79	0
43	l6	157/175 (89%)	0.19	5 (3%)	51	44	49, 57, 75, 85	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
44	L7	222/243 (91%)	0.19	3 (1%) 78 73	40, 49, 73, 106	0
44	l7	223/243 (91%)	-0.05	1 (0%) 93 92	39, 49, 80, 107	0
45	L8	233/255 (91%)	0.94	41 (17%) 2 1	67, 81, 104, 112	0
45	l8	231/255 (90%)	1.04	41 (17%) 2 1	73, 84, 106, 114	0
46	L9	191/191 (100%)	0.21	9 (4%) 35 29	60, 70, 82, 93	0
46	l9	191/191 (100%)	0.08	1 (0%) 91 90	44, 55, 71, 81	0
47	M0	211/220 (95%)	1.09	40 (18%) 2 1	45, 63, 92, 106	0
47	m0	213/220 (96%)	0.96	28 (13%) 5 3	40, 55, 80, 90	0
48	M1	169/173 (97%)	1.03	25 (14%) 3 3	68, 84, 94, 98	0
48	m1	169/173 (97%)	0.38	6 (3%) 46 39	52, 67, 77, 82	0
49	M3	193/198 (97%)	0.40	8 (4%) 41 34	39, 64, 92, 113	0
49	m3	194/198 (97%)	0.77	29 (14%) 3 2	41, 69, 100, 111	0
50	M4	136/137 (99%)	-0.04	2 (1%) 76 71	54, 61, 73, 83	0
50	m4	137/137 (100%)	-0.03	1 (0%) 89 86	48, 54, 69, 82	0
51	M5	203/203 (100%)	0.27	3 (1%) 76 71	41, 56, 66, 69	0
51	m5	203/203 (100%)	0.47	10 (4%) 33 27	44, 60, 71, 76	0
52	M6	197/198 (99%)	0.27	4 (2%) 68 62	42, 49, 65, 68	0
52	m6	197/198 (99%)	0.08	2 (1%) 84 80	34, 42, 64, 66	0
53	M7	183/183 (100%)	0.87	23 (12%) 5 4	47, 55, 98, 116	0
53	m7	155/183 (84%)	0.17	0 100 100	39, 46, 59, 77	0
54	M8	185/185 (100%)	0.22	3 (1%) 74 69	43, 55, 69, 88	0
54	m8	185/185 (100%)	0.11	0 100 100	42, 57, 66, 70	0
55	M9	188/188 (100%)	0.55	19 (10%) 9 8	65, 77, 135, 140	0
55	m9	188/188 (100%)	0.49	12 (6%) 23 19	51, 65, 121, 134	0
56	N0	172/172 (100%)	0.18	4 (2%) 64 57	49, 57, 69, 74	0
56	n0	172/172 (100%)	0.02	1 (0%) 90 88	43, 49, 59, 66	0
57	N1	159/159 (100%)	0.44	4 (2%) 61 54	42, 56, 93, 101	0
57	n1	159/159 (100%)	0.57	11 (6%) 20 16	40, 49, 83, 87	0
58	N2	100/120 (83%)	1.06	25 (25%) 1 1	93, 104, 120, 122	0
58	n2	98/120 (81%)	1.04	21 (21%) 1 1	75, 88, 94, 97	0
59	N3	136/136 (100%)	0.67	11 (8%) 15 11	51, 60, 70, 74	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
59	n3	136/136 (100%)	0.41	3 (2%) 65 59	36, 45, 55, 58	0
60	N4	98/155 (63%)	1.83	28 (28%) 1 1	62, 72, 134, 147	0
60	n4	135/155 (87%)	0.74	19 (14%) 4 3	46, 91, 115, 116	0
61	N5	121/141 (85%)	0.80	13 (10%) 8 6	58, 69, 84, 95	0
61	n5	120/141 (85%)	0.27	1 (0%) 87 84	56, 68, 85, 93	0
62	N6	126/126 (100%)	0.84	19 (15%) 3 2	50, 62, 74, 79	0
62	n6	126/126 (100%)	0.42	4 (3%) 51 44	51, 64, 77, 83	0
63	N7	135/135 (100%)	1.78	49 (36%) 0 1	82, 94, 103, 109	0
63	n7	135/135 (100%)	1.44	43 (31%) 1 1	76, 91, 102, 105	0
64	N8	148/148 (100%)	0.36	2 (1%) 78 73	35, 56, 74, 82	0
64	n8	148/148 (100%)	0.16	3 (2%) 68 62	34, 59, 74, 77	0
65	N9	58/58 (100%)	0.54	4 (6%) 20 16	38, 61, 95, 106	0
65	n9	58/58 (100%)	0.36	5 (8%) 13 10	38, 58, 78, 85	0
66	O0	97/104 (93%)	0.64	9 (9%) 11 9	81, 89, 101, 102	0
66	o0	100/104 (96%)	0.44	2 (2%) 68 62	75, 82, 102, 108	0
67	O1	109/112 (97%)	0.97	14 (12%) 5 4	61, 72, 91, 103	0
67	o1	109/112 (97%)	0.45	4 (3%) 45 38	47, 58, 81, 90	0
68	O2	127/129 (98%)	0.16	2 (1%) 74 69	36, 49, 62, 68	0
68	o2	127/129 (98%)	0.54	7 (5%) 29 23	35, 52, 63, 71	0
69	O3	106/106 (100%)	0.31	1 (0%) 85 82	42, 47, 71, 78	0
69	o3	106/106 (100%)	0.14	0 100 100	40, 46, 69, 75	0
70	O4	112/120 (93%)	0.60	7 (6%) 23 19	57, 75, 101, 107	0
70	o4	112/120 (93%)	0.66	8 (7%) 19 15	50, 68, 99, 103	0
71	O5	119/119 (100%)	0.32	6 (5%) 32 26	55, 70, 78, 82	0
71	o5	119/119 (100%)	0.59	8 (6%) 21 17	59, 71, 87, 96	0
72	O6	99/99 (100%)	0.63	5 (5%) 32 25	60, 68, 91, 102	0
72	o6	99/99 (100%)	0.40	4 (4%) 42 34	64, 72, 88, 103	0
73	O7	87/87 (100%)	0.13	1 (1%) 82 78	43, 50, 65, 71	0
73	o7	87/87 (100%)	0.21	2 (2%) 64 57	41, 50, 73, 84	0
74	O8	77/77 (100%)	0.86	8 (10%) 8 7	83, 93, 108, 113	0
74	o8	77/77 (100%)	1.51	25 (32%) 1 1	77, 86, 94, 98	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
75	O9	50/50 (100%)	0.49	1 (2%) 68 62	52, 57, 59, 60	0
75	o9	50/50 (100%)	0.27	1 (2%) 68 62	50, 55, 63, 63	0
76	Q0	52/52 (100%)	0.86	7 (13%) 4 3	53, 59, 74, 79	0
76	q0	52/52 (100%)	0.37	1 (1%) 70 63	41, 45, 55, 58	0
77	Q1	25/25 (100%)	1.26	6 (24%) 1 1	62, 66, 73, 74	0
77	q1	25/25 (100%)	0.99	2 (8%) 15 12	50, 56, 58, 58	0
78	Q2	105/105 (100%)	0.12	2 (1%) 70 63	41, 56, 76, 91	0
78	q2	105/105 (100%)	0.26	1 (0%) 84 80	44, 56, 71, 93	0
79	Q3	91/91 (100%)	0.21	2 (2%) 65 59	54, 65, 80, 85	0
79	q3	91/91 (100%)	0.38	3 (3%) 50 43	46, 59, 72, 84	0
80	m2	0/165	-	-	-	-
81	p0	120/311 (38%)	1.26	32 (26%) 1 1	101, 124, 144, 152	0
82	p1	0/47	-	-	-	-
82	p2	0/47	-	-	-	-
83	f	148/157 (94%)	2.86	92 (62%) 0 0	55, 96, 150, 153	75 (50%)
84	B	2/5 (40%)	0.21	0 100 100	43, 43, 43, 44	0
84	C	2/5 (40%)	0.25	0 100 100	49, 49, 49, 49	0
All	All	33084/35503 (93%)	0.60	3317 (10%) 9 8	34, 74, 132, 212	75 (0%)

All (3317) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
60	N4	75	THR	13.9
1	2	1699	G	13.8
53	M7	161	ALA	12.0
1	2	1696	G	12.0
1	2	1697	G	11.7
60	N4	86	SER	11.4
53	M7	160	ALA	11.3
60	N4	88	ASP	11.0
1	2	1709	C	11.0
71	o5	120	ALA	10.8
32	E0	61	SER	10.8
36	1	1570	U	10.5
1	2	1698	G	10.4
12	c0	23	ALA	10.2
34	sR	213	SER	9.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
60	N4	89	LEU	9.7
1	2	1708	U	9.7
36	5	2443	A	9.6
36	5	2442	G	9.5
33	e1	77	GLY	9.4
7	S5	152	GLY	9.4
53	M7	184	ALA	9.2
1	2	1694	A	9.2
16	C4	75	GLY	9.1
32	e0	63	GLN	9.1
83	f	120	LEU	9.0
13	c1	3	THR	9.0
34	SR	32	LEU	8.9
7	S5	36	ALA	8.7
13	c1	4	GLU	8.6
18	C6	20	ALA	8.4
53	M7	162	GLU	8.3
1	2	1702	A	8.3
31	d9	29	GLY	8.2
1	6	662	U	8.2
34	SR	25	THR	8.2
36	1	1568	U	8.1
34	sR	314	GLN	8.0
60	N4	83	THR	7.9
18	C6	21	HIS	7.9
7	S5	153	GLY	7.7
34	SR	44	SER	7.7
83	f	134	LEU	7.7
3	s1	89	ASP	7.6
1	2	1695	G	7.6
21	c9	55	TYR	7.6
60	N4	87	LEU	7.4
60	N4	85	ALA	7.4
13	c1	2	SER	7.3
27	d5	50	ILE	7.3
3	S1	20	VAL	7.2
22	d0	67	THR	7.2
12	c0	22	VAL	7.2
34	SR	262	VAL	7.1
28	D6	20	PRO	7.1
1	6	678	A	7.1
28	d6	63	ALA	7.0

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Mol	Chain	Res	Type	RSRZ
22	d0	14	GLN	7.0
34	sR	24	ALA	6.9
36	5	2494	A	6.9
21	c9	18	TYR	6.8
1	6	659	C	6.8
42	L5	6	ASP	6.8
7	S5	71	ALA	6.8
1	2	1700	C	6.8
6	S4	54	TYR	6.8
22	D0	120	SER	6.7
12	c0	65	TYR	6.7
19	C7	99	VAL	6.6
11	S9	87	SER	6.6
83	f	31	PHE	6.6
60	N4	97	LYS	6.6
17	c5	134	THR	6.6
83	f	121	GLY	6.5
28	D6	76	SER	6.5
60	N4	81	PRO	6.5
22	d0	98	GLN	6.5
34	SR	131	ILE	6.4
22	d0	18	GLN	6.4
35	sM	85	SER	6.4
22	d0	100	VAL	6.3
32	E0	47	VAL	6.3
7	s5	152	GLY	6.3
32	E0	60	PRO	6.3
83	f	71	GLU	6.3
29	d7	59	CYS	6.2
34	sR	303	ALA	6.3
45	l8	192	GLN	6.2
60	N4	76	VAL	6.2
1	2	913	G	6.2
7	S5	151	GLY	6.2
36	5	2444	C	6.2
22	d0	101	LYS	6.2
29	D7	38	PRO	6.1
60	N4	72	SER	6.1
60	N4	74	LYS	6.1
1	2	1701	A	6.1
36	5	2496	C	6.1
7	S5	37	GLN	6.1

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Mol	Chain	Res	Type	RSRZ
3	s1	54	LEU	6.1
83	f	30	GLY	6.1
83	f	38	PRO	6.1
36	5	2495	C	6.1
1	6	663	U	6.1
83	f	86	LYS	6.1
29	d7	58	SER	6.0
17	C5	50	THR	6.0
6	S4	111	VAL	6.0
36	1	1569	U	6.0
34	sR	302	PHE	6.0
13	C1	2	SER	6.0
22	d0	95	ALA	6.0
39	L2	253	GLN	6.0
1	2	506	A	6.0
6	S4	77	ARG	6.0
4	S2	62	PRO	6.0
83	f	83	PRO	6.0
18	C6	8	GLN	6.0
36	5	2441	A	5.9
45	l8	120	LYS	5.9
1	2	1703	C	5.9
83	f	22	GLN	5.9
60	N4	82	ILE	5.9
16	C4	15	GLY	5.9
35	sM	40	PRO	5.9
22	d0	93	LEU	5.9
6	S4	66	MET	5.9
16	c4	55	SER	5.9
32	E0	46	ASN	5.9
6	S4	71	LYS	5.8
83	f	64	ILE	5.8
3	S1	94	LYS	5.8
22	d0	64	LYS	5.8
12	c0	20	VAL	5.8
63	n7	2	ALA	5.8
48	M1	127	PHE	5.8
12	c0	25	LYS	5.8
81	p0	69	ASP	5.8
16	C4	16	VAL	5.8
42	L5	51	LEU	5.8
83	f	157	ASP	5.7

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
31	d9	30	LEU	5.7
6	s4	15	PRO	5.7
4	s2	105	GLY	5.7
23	d1	87	ARG	5.7
83	f	135	MET	5.7
83	f	149	SER	5.6
34	sR	214	ALA	5.6
1	6	1491	U	5.6
19	C7	121	VAL	5.6
1	2	1705	C	5.6
1	2	1707	A	5.6
1	2	1706	C	5.6
83	f	90	TYR	5.6
5	s3	37	VAL	5.6
19	C7	62	GLN	5.6
5	s3	69	LEU	5.5
36	1	1567	U	5.5
81	p0	88	PHE	5.5
16	c4	28	VAL	5.5
29	D7	37	CYS	5.5
16	C4	74	VAL	5.5
8	S6	175	ILE	5.5
11	s9	148	VAL	5.5
36	5	2447	A	5.5
34	sR	158	PRO	5.5
45	L8	116	VAL	5.5
81	p0	68	SER	5.5
33	e1	112	GLY	5.5
6	S4	65	LEU	5.5
1	6	493	U	5.4
6	s4	261	LEU	5.4
22	d0	19	ILE	5.4
11	S9	128	LEU	5.4
9	s7	43	PHE	5.4
34	sR	170	ILE	5.4
18	C6	143	ARG	5.4
12	c0	57	THR	5.3
53	M7	168	LEU	5.3
2	S0	98	ILE	5.3
1	2	1690	G	5.3
3	S1	217	LEU	5.3
36	5	2451	G	5.3

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Mol	Chain	Res	Type	RSRZ
11	S9	85	VAL	5.3
19	C7	2	GLY	5.3
2	S0	26	ALA	5.3
8	S6	149	LYS	5.3
3	S1	50	LYS	5.3
1	2	1693	A	5.3
3	S1	140	ILE	5.3
6	s4	131	LEU	5.3
22	D0	121	ASN	5.3
35	SM	86	ASN	5.3
7	S5	96	SER	5.3
9	s7	70	PHE	5.3
28	D6	49	ALA	5.3
60	n4	69	LYS	5.3
83	f	87	ARG	5.3
1	6	658	C	5.3
2	s0	20	ALA	5.3
20	C8	2	SER	5.2
83	f	136	VAL	5.2
7	s5	150	GLY	5.2
28	D6	31	PRO	5.2
12	c0	67	THR	5.2
34	SR	33	LEU	5.2
83	f	142	MET	5.2
31	d9	20	GLN	5.2
34	sR	244	ALA	5.2
56	N0	1	MET	5.2
44	L7	28	ALA	5.2
18	C6	22	VAL	5.2
63	N7	26	VAL	5.2
83	f	18	THR	5.2
10	S8	168	CYS	5.2
8	S6	77	LEU	5.2
7	S5	70	VAL	5.2
45	l8	109	LEU	5.2
34	sR	202	LEU	5.1
6	S4	102	VAL	5.1
1	6	495	C	5.1
35	sM	39	PRO	5.1
27	d5	89	ILE	5.1
1	2	1711	C	5.1
14	c2	65	SER	5.1

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Mol	Chain	Res	Type	RSRZ
11	S9	86	LEU	5.1
83	f	81	GLU	5.1
5	s3	175	VAL	5.1
13	c1	5	LEU	5.1
29	d7	46	VAL	5.1
10	s8	176	SER	5.1
30	D8	67	ARG	5.1
4	S2	63	VAL	5.1
3	S1	55	LYS	5.0
81	p0	20	GLU	5.0
27	d5	51	LEU	5.0
83	f	36	SER	5.0
34	sR	72	THR	5.0
1	2	1692	G	5.0
20	C8	146	ALA	5.0
35	SM	89	ARG	5.0
83	f	41	ILE	5.0
2	S0	100	GLY	5.0
12	C0	22	VAL	5.0
45	l8	121	SER	5.0
34	sR	252	LEU	5.0
36	5	2452	G	5.0
12	c0	68	LEU	5.0
32	E0	56	MET	5.0
8	S6	73	ILE	5.0
10	s8	200	LYS	5.0
20	C8	15	LEU	5.0
36	5	2440	G	5.0
6	S4	261	LEU	5.0
22	D0	82	TYR	5.0
18	C6	15	SER	5.0
83	f	23	CYS	5.0
22	d0	97	VAL	5.0
22	d0	96	PRO	5.0
36	5	1569	U	5.0
23	D1	56	SER	4.9
4	S2	232	GLU	4.9
5	s3	6	SER	4.9
10	s8	143	TRP	4.9
6	S4	69	HIS	4.9
22	d0	54	GLY	4.9
34	SR	61	PHE	4.9

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Mol	Chain	Res	Type	RSRZ
83	f	153	ALA	4.9
34	sR	166	SER	4.9
9	s7	69	GLY	4.9
42	L5	3	PHE	4.9
18	C6	68	ARG	4.9
2	s0	98	ILE	4.9
27	d5	42	LEU	4.9
12	C0	24	LYS	4.9
74	O8	43	PHE	4.9
83	f	72	ASP	4.9
83	f	42	VAL	4.9
2	s0	170	ILE	4.9
21	c9	28	LEU	4.9
42	L5	146	LEU	4.9
34	sR	313	TRP	4.9
49	m3	131	LYS	4.9
63	N7	2	ALA	4.9
8	S6	75	LEU	4.9
60	n4	67	VAL	4.9
2	s0	46	HIS	4.9
35	SM	85	SER	4.9
11	S9	186	GLU	4.9
11	S9	96	VAL	4.8
22	d0	121	ASN	4.8
1	2	1710	U	4.8
33	e1	79	LYS	4.8
3	S1	86	LEU	4.8
81	p0	70	LEU	4.8
11	S9	2	PRO	4.8
1	2	277	U	4.8
3	S1	233	GLY	4.8
31	D9	4	GLU	4.8
45	L8	202	GLU	4.8
34	sR	212	ALA	4.8
30	d8	9	LEU	4.8
31	d9	27	HIS	4.8
6	S4	101	LEU	4.8
34	sR	157	VAL	4.8
36	5	1567	U	4.8
3	S1	220	GLN	4.8
17	C5	28	MET	4.8
12	C0	64	TYR	4.8

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Mol	Chain	Res	Type	RSRZ
21	c9	118	PRO	4.8
22	d0	102	ARG	4.7
71	o5	119	LYS	4.7
10	s8	177	GLY	4.7
53	M7	183	ALA	4.7
6	S4	47	PHE	4.7
28	d6	69	ASN	4.7
30	d8	44	VAL	4.7
34	SR	253	ALA	4.7
26	D4	26	ASP	4.7
58	n2	14	THR	4.7
45	L8	93	LEU	4.7
21	c9	25	GLN	4.7
16	c4	105	LEU	4.7
23	D1	55	LEU	4.7
6	s4	14	ALA	4.7
26	D4	18	LEU	4.7
34	sR	301	LEU	4.7
14	c2	123	VAL	4.7
17	c5	135	THR	4.7
5	s3	38	GLU	4.7
34	SR	79	TYR	4.7
83	f	85	VAL	4.7
34	SR	115	ILE	4.7
16	C4	34	SER	4.7
8	S6	148	SER	4.7
29	D7	48	SER	4.6
28	d6	44	ILE	4.6
45	l8	111	LYS	4.6
47	m0	173	PHE	4.6
5	s3	184	ILE	4.6
2	S0	97	PRO	4.6
83	f	32	VAL	4.6
7	S5	154	ALA	4.6
83	f	105	MET	4.6
34	SR	55	GLY	4.6
3	S1	92	GLN	4.6
34	sR	26	SER	4.6
11	S9	156	ILE	4.6
45	L8	26	LEU	4.6
74	o8	2	ALA	4.6
10	S8	152	ILE	4.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
22	d0	78	THR	4.6
49	m3	130	GLY	4.6
49	m3	133	PRO	4.6
5	s3	11	LEU	4.6
34	sR	11	GLY	4.6
83	f	65	PHE	4.6
1	2	715	U	4.6
5	s3	43	PRO	4.6
22	D0	86	ILE	4.6
42	L5	148	ILE	4.5
3	S1	138	PHE	4.5
28	D6	9	GLY	4.5
16	C4	29	HIS	4.5
1	2	820	U	4.5
6	s4	183	VAL	4.5
81	p0	19	LEU	4.5
28	D6	17	HIS	4.5
63	N7	124	ALA	4.5
2	S0	18	LEU	4.5
60	N4	70	LYS	4.5
18	c6	44	LEU	4.5
22	d0	27	THR	4.5
3	S1	151	LYS	4.5
21	C9	28	LEU	4.5
34	sR	210	LEU	4.5
63	n7	118	PHE	4.5
12	C0	6	GLU	4.5
26	D4	25	VAL	4.5
12	c0	64	TYR	4.5
28	D6	21	VAL	4.5
5	s3	42	THR	4.5
1	2	709	C	4.5
71	o5	118	ILE	4.5
23	D1	53	TYR	4.5
2	S0	201	LEU	4.4
81	p0	86	PHE	4.4
43	L6	131	LYS	4.4
21	C9	108	LEU	4.4
5	s3	113	LEU	4.4
60	N4	90	ILE	4.4
63	n7	75	VAL	4.4
36	1	1566	A	4.4

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Mol	Chain	Res	Type	RSRZ
8	S6	78	THR	4.4
47	M0	33	ILE	4.4
53	M7	176	ILE	4.4
36	1	1573	G	4.4
34	sR	13	LEU	4.4
47	m0	195	ALA	4.4
83	f	115	ALA	4.4
10	s8	67	TRP	4.4
83	f	43	ASP	4.4
3	S1	110	LEU	4.4
83	f	147	ALA	4.4
2	s0	22	THR	4.4
10	s8	69	SER	4.4
45	L8	94	PHE	4.4
9	S7	150	GLN	4.4
22	d0	28	SER	4.4
63	N7	46	ILE	4.4
5	S3	21	LEU	4.4
36	1	1565	G	4.4
60	N4	96	LEU	4.4
36	5	2446	U	4.3
5	s3	134	CYS	4.3
83	f	70	LEU	4.3
4	s2	88	LYS	4.3
5	s3	8	LYS	4.3
8	S6	74	LYS	4.3
31	d9	16	LYS	4.3
9	s7	54	GLY	4.3
16	C4	115	ILE	4.3
36	5	2539	C	4.3
11	S9	106	GLU	4.3
14	c2	128	ALA	4.3
7	s5	37	GLN	4.3
18	C6	7	VAL	4.3
30	D8	16	LEU	4.3
45	l8	203	VAL	4.3
63	n7	68	ILE	4.3
31	d9	14	TYR	4.3
78	Q2	36	PHE	4.3
4	S2	145	GLY	4.3
11	s9	156	ILE	4.3
53	M7	167	ARG	4.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
7	S5	41	LYS	4.3
12	c0	26	ASP	4.3
1	2	1794	A	4.3
81	p0	81	LYS	4.3
34	SR	278	PHE	4.3
22	d0	20	ILE	4.3
29	D7	43	ILE	4.3
63	N7	91	ALA	4.3
76	Q0	77	ILE	4.3
19	C7	123	ASN	4.3
3	S1	28	GLU	4.3
83	f	39	CYS	4.3
1	2	1691	A	4.3
6	S4	78	THR	4.3
35	SM	87	THR	4.3
83	f	156	THR	4.3
28	D6	69	ASN	4.3
1	2	134	U	4.3
18	C6	57	LEU	4.3
1	2	276	C	4.3
12	C0	66	TYR	4.3
6	S4	52	LEU	4.2
48	M1	96	PHE	4.2
61	N5	107	VAL	4.2
63	N7	113	VAL	4.2
5	s3	98	ALA	4.2
83	f	11	ALA	4.2
31	d9	40	ARG	4.2
63	n7	96	VAL	4.2
32	E0	45	VAL	4.2
8	s6	215	ARG	4.2
74	O8	6	THR	4.2
59	N3	2	SER	4.2
36	5	1350	A	4.2
40	l3	387	LEU	4.2
36	5	2493	U	4.2
3	s1	51	SER	4.2
63	N7	99	GLU	4.2
22	d0	87	HIS	4.2
76	Q0	128	LYS	4.2
31	d9	36	LEU	4.2
67	O1	71	LEU	4.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
36	5	2445	A	4.2
21	c9	22	LEU	4.2
34	SR	198	ASN	4.2
63	N7	92	PHE	4.2
36	5	1352	A	4.2
83	f	138	ILE	4.2
2	S0	198	MET	4.2
34	SR	78	ALA	4.2
33	e1	88	PRO	4.2
45	l8	200	LEU	4.2
16	C4	18	ARG	4.2
28	D6	18	VAL	4.2
63	N7	65	ARG	4.2
39	l2	164	GLY	4.2
3	S1	47	LEU	4.2
34	SR	45	TRP	4.2
60	N4	98	PRO	4.2
12	C0	5	LYS	4.2
11	S9	148	VAL	4.2
12	C0	16	PHE	4.2
34	SR	46	LYS	4.2
36	5	2498	U	4.2
11	S9	99	LEU	4.2
18	C6	92	TYR	4.2
34	SR	94	VAL	4.2
42	L5	101	THR	4.2
42	L5	92	LEU	4.1
9	S7	108	GLN	4.1
2	s0	165	ARG	4.1
4	S2	45	VAL	4.1
33	E1	91	ILE	4.1
6	S4	63	ALA	4.1
34	sR	272	ASP	4.1
1	6	718	U	4.1
43	L6	130	ILE	4.1
32	E0	48	THR	4.1
63	n7	92	PHE	4.1
83	f	28	LYS	4.1
36	5	1568	U	4.1
21	C9	6	VAL	4.1
49	m3	95	ILE	4.1
6	S4	67	GLN	4.1

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Mol	Chain	Res	Type	RSRZ
31	d9	52	PHE	4.1
44	L7	25	GLN	4.1
45	l8	94	PHE	4.1
63	N7	120	GLU	4.1
83	f	45	SER	4.1
12	c0	62	GLN	4.1
3	S1	114	VAL	4.1
3	S1	89	ASP	4.1
6	S4	49	ARG	4.1
6	S4	79	ASP	4.1
26	D4	16	PRO	4.1
21	c9	37	VAL	4.1
63	N7	7	ALA	4.1
71	O5	120	ALA	4.1
55	m9	184	LEU	4.1
63	N7	5	LEU	4.1
3	S1	52	THR	4.1
48	M1	104	PHE	4.1
8	S6	186	ARG	4.1
10	s8	179	CYS	4.1
50	M4	138	ALA	4.1
83	f	16	SER	4.1
21	c9	122	ARG	4.1
27	D5	58	ARG	4.1
35	sM	46	LYS	4.1
52	m6	182	ASN	4.1
58	N2	89	LEU	4.1
83	f	46	THR	4.0
5	S3	186	VAL	4.0
63	N7	132	SER	4.0
79	q3	2	ALA	4.0
14	C2	32	LEU	4.0
20	C8	145	ARG	4.0
13	C1	4	GLU	4.0
36	5	1016	C	4.0
20	C8	22	VAL	4.0
1	2	914	G	4.0
3	s1	84	ILE	4.0
9	S7	98	ILE	4.0
12	C0	62	GLN	4.0
27	D5	89	ILE	4.0
32	E0	32	GLY	4.0

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Mol	Chain	Res	Type	RSRZ
1	2	910	C	4.0
29	d7	24	LEU	4.0
42	L5	95	TRP	4.0
45	l8	104	GLU	4.0
74	o8	34	ALA	4.0
11	S9	138	LYS	4.0
18	C6	114	ARG	4.0
35	SM	84	LYS	4.0
45	l8	142	LEU	4.0
16	c4	27	PHE	4.0
18	c6	141	SER	4.0
7	S5	150	GLY	4.0
2	s0	48	ILE	4.0
6	S4	44	LEU	4.0
11	S9	144	PRO	4.0
5	s3	9	ARG	4.0
11	s9	111	THR	4.0
14	c2	103	LEU	4.0
45	L8	131	ALA	4.0
5	s3	79	TYR	4.0
74	o8	33	LYS	4.0
83	f	119	GLU	4.0
55	M9	51	VAL	4.0
6	S4	26	CYS	4.0
1	6	664	U	4.0
3	S1	101	HIS	4.0
11	S9	29	LYS	4.0
18	c6	142	TYR	4.0
63	n7	49	TYR	4.0
1	6	1694	A	4.0
20	c8	22	VAL	4.0
5	s3	86	LEU	4.0
3	S1	139	ALA	4.0
5	S3	24	PHE	4.0
21	c9	111	ILE	4.0
19	C7	65	PRO	4.0
35	sM	38	PRO	4.0
3	S1	141	ALA	4.0
5	s3	65	ARG	4.0
16	C4	119	THR	4.0
42	L5	60	ILE	4.0
8	S6	199	GLN	3.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	s1	217	LEU	3.9
16	c4	53	ASP	3.9
81	p0	192	ASP	3.9
25	D3	48	HIS	3.9
36	5	2540	A	3.9
11	S9	141	VAL	3.9
83	f	73	LEU	3.9
34	SR	72	THR	3.9
9	S7	115	SER	3.9
33	E1	82	LYS	3.9
36	5	1566	A	3.9
74	o8	54	LEU	3.9
3	S1	103	MET	3.9
21	c9	21	PHE	3.9
3	S1	100	PHE	3.9
19	C7	71	PHE	3.9
2	S0	99	ALA	3.9
31	d9	28	THR	3.9
11	S9	135	ALA	3.9
2	S0	158	VAL	3.9
18	C6	79	TYR	3.9
11	S9	28	LEU	3.9
21	c9	4	VAL	3.9
16	C4	97	GLY	3.9
3	s1	47	LEU	3.9
9	s7	41	LEU	3.9
26	D4	35	VAL	3.9
39	l2	253	GLN	3.9
17	c5	85	ILE	3.9
26	D4	7	ILE	3.9
83	f	20	PRO	3.9
11	S9	97	LEU	3.9
83	f	84	VAL	3.9
22	D0	54	GLY	3.9
34	SR	263	PHE	3.9
2	s0	162	CYS	3.9
13	C1	145	ALA	3.9
32	E0	25	GLU	3.9
36	1	2502	A	3.9
28	D6	80	HIS	3.9
6	S4	45	ILE	3.9
22	d0	103	ILE	3.9

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Mol	Chain	Res	Type	RSRZ
53	M7	182	ILE	3.9
40	L3	337	THR	3.9
29	D7	41	LEU	3.9
60	n4	134	GLN	3.9
34	SR	284	ALA	3.9
42	L5	144	VAL	3.9
28	D6	2	PRO	3.9
6	s4	162	ILE	3.9
36	1	1571	A	3.9
55	M9	178	ALA	3.9
83	f	19	TYR	3.8
8	s6	201	GLN	3.8
11	S9	180	LYS	3.8
16	C4	46	MET	3.8
1	6	1199	G	3.8
13	C1	3	THR	3.8
52	M6	184	THR	3.8
53	M7	181	ARG	3.8
3	S1	91	VAL	3.8
32	E0	44	PHE	3.8
21	C9	27	LYS	3.8
27	D5	88	ILE	3.8
16	c4	42	VAL	3.8
7	S5	149	VAL	3.8
70	O4	73	SER	3.8
34	SR	214	ALA	3.8
40	L3	217	ALA	3.8
35	sM	33	LYS	3.8
26	D4	31	ASN	3.8
30	D8	66	LEU	3.8
34	SR	313	TRP	3.8
1	2	1059	U	3.8
35	sM	28	SER	3.8
21	C9	95	ASP	3.8
20	c8	18	LEU	3.8
36	1	1349	G	3.8
47	M0	87	LEU	3.8
5	s3	85	VAL	3.8
83	f	35	LYS	3.8
7	S5	155	ALA	3.8
15	C3	15	ALA	3.8
45	l8	110	THR	3.8

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Mol	Chain	Res	Type	RSRZ
60	n4	68	ALA	3.8
36	1	1572	U	3.8
63	N7	42	LEU	3.8
3	S1	142	PHE	3.8
3	s1	30	PHE	3.8
34	sR	253	ALA	3.8
65	N9	55	ALA	3.8
71	O5	20	GLN	3.8
40	L3	164	THR	3.8
45	L8	126	SER	3.8
2	S0	122	ILE	3.8
83	f	97	ASP	3.8
34	sR	315	VAL	3.8
1	6	1700	C	3.8
6	S4	55	ALA	3.8
11	S9	115	LYS	3.8
28	D6	67	THR	3.8
28	D6	89	ARG	3.8
28	d6	64	LEU	3.8
26	D4	72	PHE	3.8
28	D6	8	ASN	3.8
34	sR	229	LYS	3.8
22	d0	107	THR	3.8
27	d5	102	THR	3.8
5	s3	75	LYS	3.8
5	s3	179	GLN	3.7
45	L8	130	TYR	3.7
18	C6	65	ILE	3.7
6	S4	80	THR	3.7
33	E1	87	THR	3.7
6	S4	25	GLY	3.7
34	sR	55	GLY	3.7
12	C0	13	GLN	3.7
34	sR	211	ILE	3.7
66	O0	62	LEU	3.7
45	L8	91	PHE	3.7
47	M0	148	VAL	3.7
63	n7	130	PHE	3.7
42	L5	89	THR	3.7
7	s5	159	ALA	3.7
7	s5	158	GLN	3.7
7	S5	222	LYS	3.7

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
12	C0	76	LEU	3.7
18	C6	36	ILE	3.7
49	m3	54	LEU	3.7
83	f	139	ILE	3.7
83	f	57	VAL	3.7
6	S4	64	ILE	3.7
12	c0	66	TYR	3.7
6	s4	139	VAL	3.7
48	M1	148	VAL	3.7
53	M7	157	VAL	3.7
3	S1	95	ASN	3.7
45	L8	191	ASN	3.7
34	SR	121	MET	3.7
60	n4	132	GLY	3.7
21	c9	93	HIS	3.7
58	n2	13	LYS	3.7
5	s3	151	LYS	3.7
11	S9	167	ALA	3.7
74	o8	53	THR	3.7
5	s3	138	VAL	3.7
22	D0	71	PRO	3.7
1	2	719	U	3.7
31	D9	5	ASN	3.7
36	1	1252	A	3.7
2	S0	17	LEU	3.7
6	S4	180	LEU	3.7
70	o4	38	LEU	3.7
7	S5	69	PHE	3.7
18	C6	9	THR	3.7
18	c6	90	VAL	3.7
19	c7	25	THR	3.7
83	f	63	ASP	3.7
83	f	44	MET	3.7
34	sR	81	LEU	3.7
58	n2	17	VAL	3.7
9	S7	31	SER	3.7
22	d0	13	GLU	3.7
26	D4	69	SER	3.7
32	E0	30	PRO	3.7
34	SR	92	TRP	3.7
43	l6	129	GLU	3.7
5	s3	114	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
15	C3	57	ALA	3.7
10	s8	60	ILE	3.7
67	O1	14	ILE	3.7
83	f	148	ILE	3.7
8	S6	67	VAL	3.7
45	L8	197	VAL	3.7
2	s0	101	ARG	3.7
7	s5	151	GLY	3.7
5	s3	182	LEU	3.7
18	c6	52	LEU	3.7
11	s9	91	LYS	3.7
22	d0	99	ILE	3.7
34	sR	115	ILE	3.7
83	f	33	VAL	3.7
11	s9	112	GLN	3.7
6	s4	132	GLY	3.7
3	S1	98	THR	3.7
17	C5	52	LYS	3.7
83	f	17	ALA	3.7
26	D4	70	VAL	3.7
48	M1	167	TYR	3.7
12	c0	28	ASN	3.7
42	L5	151	GLN	3.7
34	SR	7	LEU	3.6
21	C9	21	PHE	3.6
22	d0	22	ILE	3.6
26	D4	82	ALA	3.6
10	s8	103	GLN	3.6
19	C7	22	PRO	3.6
55	m9	178	ALA	3.6
83	f	141	ALA	3.6
26	D4	57	VAL	3.6
8	S6	95	LYS	3.6
34	sR	32	LEU	3.6
49	m3	129	ASN	3.6
81	p0	26	PHE	3.6
7	S5	147	THR	3.6
21	C9	71	VAL	3.6
22	D0	81	THR	3.6
34	SR	26	SER	3.6
34	SR	246	SER	3.6
4	S2	66	PHE	3.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
27	D5	50	ILE	3.6
28	D6	70	LYS	3.6
34	SR	211	ILE	3.6
43	l6	130	ILE	3.6
58	n2	106	ALA	3.6
63	N7	133	LYS	3.6
34	SR	291	SER	3.6
12	C0	29	GLN	3.6
42	L5	83	LEU	3.6
1	2	714	G	3.6
10	S8	179	CYS	3.6
3	s1	91	VAL	3.6
63	n7	95	VAL	3.6
34	sR	294	TRP	3.6
2	S0	188	LEU	3.6
22	D0	70	THR	3.6
42	L5	103	LEU	3.6
28	D6	83	ILE	3.6
27	D5	102	THR	3.6
29	D7	45	THR	3.6
8	s6	162	VAL	3.6
28	D6	73	TYR	3.6
28	D6	82	ARG	3.6
15	c3	59	GLY	3.6
17	c5	125	PRO	3.6
83	f	59	LEU	3.6
11	S9	112	GLN	3.6
5	s3	84	ILE	3.6
6	S4	162	ILE	3.6
6	S4	76	VAL	3.6
21	C9	39	THR	3.6
23	d1	39	VAL	3.6
39	l2	252	THR	3.6
42	L5	86	TYR	3.6
1	6	660	G	3.6
36	1	1243	G	3.6
3	S1	223	PHE	3.6
19	C7	83	GLN	3.6
27	D5	103	ARG	3.6
23	D1	34	ILE	3.6
33	e1	86	THR	3.6
57	n1	33	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
61	N5	124	VAL	3.6
11	S9	181	ALA	3.6
12	c0	24	LYS	3.6
21	C9	100	ILE	3.6
6	s4	13	ALA	3.6
17	C5	131	ALA	3.6
1	6	661	A	3.5
5	s3	62	ASN	3.5
3	S1	58	SER	3.5
6	S4	166	SER	3.5
28	D6	92	ARG	3.5
5	s3	77	PHE	3.5
49	m3	145	PHE	3.5
36	5	2506	U	3.5
59	N3	81	GLN	3.5
48	M1	55	ARG	3.5
6	S4	91	THR	3.5
12	C0	67	THR	3.5
1	2	718	U	3.5
19	C7	74	GLN	3.5
35	SM	90	ALA	3.5
3	s1	100	PHE	3.5
11	S9	146	PHE	3.5
49	m3	119	TYR	3.5
81	p0	24	SER	3.5
5	S3	74	GLN	3.5
3	S1	53	GLY	3.5
7	S5	165	LEU	3.5
21	C9	38	LYS	3.5
45	l8	150	LEU	3.5
36	1	1239	C	3.5
42	l5	125	VAL	3.5
29	d7	33	LEU	3.5
34	sR	243	LEU	3.5
49	m3	132	ALA	3.5
63	n7	45	GLY	3.5
83	f	13	ALA	3.5
34	SR	305	TYR	3.5
31	d9	4	GLU	3.5
6	S4	48	LEU	3.5
5	s3	152	PHE	3.5
12	C0	79	TYR	3.5

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Mol	Chain	Res	Type	RSRZ
36	1	1270	A	3.5
20	C8	40	ARG	3.5
13	c1	117	VAL	3.5
16	c4	79	VAL	3.5
70	O4	23	VAL	3.5
2	s0	146	LEU	3.5
34	sR	141	LEU	3.5
6	S4	167	GLY	3.5
34	sR	185	GLN	3.5
22	d0	116	VAL	3.5
8	s6	193	LEU	3.5
34	SR	42	LEU	3.5
47	M0	97	LEU	3.5
61	N5	108	LEU	3.5
13	C1	30	ARG	3.5
6	s4	133	LYS	3.5
34	sR	25	THR	3.5
34	sR	52	GLN	3.5
4	S2	144	TRP	3.5
20	c8	52	VAL	3.5
21	C9	124	ILE	3.5
47	M0	36	LEU	3.5
3	S1	42	ASN	3.5
12	c0	59	PHE	3.5
63	N7	45	GLY	3.5
70	o4	39	ALA	3.5
3	S1	46	THR	3.5
6	S4	15	PRO	3.5
16	C4	89	THR	3.5
33	E1	86	THR	3.5
58	n2	98	THR	3.5
11	S9	3	ARG	3.5
83	f	92	LEU	3.5
36	5	2098	C	3.4
3	s1	133	TYR	3.4
36	5	2499	U	3.4
38	4	158	U	3.4
58	N2	27	VAL	3.4
11	s9	90	LYS	3.4
14	c2	59	LEU	3.4
19	C7	63	LYS	3.4
63	n7	46	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
83	f	40	LYS	3.4
4	s2	87	GLN	3.4
14	C2	27	ALA	3.4
55	m9	189	ALA	3.4
63	n7	76	ASN	3.4
1	6	232	U	3.4
2	s0	160	ILE	3.4
20	C8	17	LEU	3.4
21	c9	92	LYS	3.4
28	D6	27	SER	3.4
34	SR	34	LEU	3.4
45	L8	90	THR	3.4
1	6	1699	G	3.4
70	O4	64	THR	3.4
6	S4	8	HIS	3.4
32	e0	2	ALA	3.4
34	SR	296	ALA	3.4
11	S9	36	LEU	3.4
12	C0	75	TYR	3.4
14	c2	121	VAL	3.4
47	M0	152	LEU	3.4
16	C4	83	ILE	3.4
28	D6	41	ILE	3.4
63	n7	131	PHE	3.4
1	2	1704	U	3.4
9	S7	48	GLU	3.4
58	n2	70	LYS	3.4
83	f	66	THR	3.4
9	S7	62	VAL	3.4
33	E1	145	HIS	3.4
51	m5	148	TYR	3.4
8	S6	66	GLY	3.4
15	C3	14	SER	3.4
1	2	1370	U	3.4
18	c6	79	TYR	3.4
26	D4	22	GLN	3.4
32	e0	17	GLN	3.4
34	SR	20	VAL	3.4
47	M0	96	VAL	3.4
11	S9	104	PHE	3.4
42	l5	247	ILE	3.4
35	SM	88	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
55	M9	52	LYS	3.4
18	C6	6	SER	3.4
81	p0	187	VAL	3.4
12	c0	58	GLN	3.4
27	d5	41	ILE	3.4
48	m1	174	LYS	3.4
58	n2	11	ILE	3.4
6	s4	101	LEU	3.4
11	s9	80	LEU	3.4
16	c4	78	ALA	3.4
20	c8	17	LEU	3.4
31	d9	23	VAL	3.4
36	5	2497	U	3.4
28	D6	90	GLU	3.4
60	n4	66	GLU	3.4
16	c4	48	VAL	3.4
21	c9	108	LEU	3.4
81	p0	87	VAL	3.4
12	C0	65	TYR	3.4
36	1	1351	U	3.4
6	s4	18	TRP	3.4
16	C4	47	LYS	3.4
28	D6	78	ALA	3.4
58	N2	92	TRP	3.4
61	N5	82	LEU	3.4
71	o5	110	ALA	3.4
7	s5	68	ILE	3.4
12	C0	35	ILE	3.4
17	c5	84	ILE	3.4
17	c5	103	ASN	3.4
19	C7	101	ASN	3.4
21	c9	100	ILE	3.4
34	sR	159	ASN	3.4
40	L3	49	TYR	3.4
2	S0	22	THR	3.4
4	s2	89	GLN	3.4
60	N4	84	GLY	3.4
34	sR	6	VAL	3.4
10	s8	78	ILE	3.4
27	d5	59	TYR	3.4
28	D6	68	TYR	3.4
34	sR	122	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
63	N7	72	ILE	3.4
16	C4	48	VAL	3.3
22	d0	84	MET	3.4
63	N7	96	VAL	3.3
74	O8	51	LEU	3.3
36	5	2500	A	3.3
42	L5	2	ALA	3.3
27	d5	88	ILE	3.3
72	o6	58	ILE	3.3
28	D6	19	LYS	3.3
33	e1	83	LYS	3.3
6	s4	26	CYS	3.3
66	o0	23	TYR	3.3
7	S5	217	LEU	3.3
9	s7	58	LEU	3.3
53	M7	169	THR	3.3
13	c1	116	ARG	3.3
19	C7	64	GLY	3.3
1	6	494	U	3.3
1	6	679	U	3.3
6	S4	57	ASN	3.3
45	l8	152	LEU	3.3
47	M0	91	VAL	3.3
34	SR	117	LYS	3.3
59	n3	2	SER	3.3
63	n7	6	LYS	3.3
74	o8	63	LYS	3.3
81	p0	73	PHE	3.3
2	S0	126	PRO	3.3
6	s4	149	TYR	3.3
63	N7	90	GLU	3.3
7	s5	161	ASP	3.3
11	S9	110	GLN	3.3
1	2	656	G	3.3
12	C0	25	LYS	3.3
22	d0	37	VAL	3.3
30	d8	61	ARG	3.3
59	N3	4	ASN	3.3
7	s5	36	ALA	3.3
3	s1	86	LEU	3.3
11	S9	80	LEU	3.3
27	d5	52	LYS	3.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
34	SR	221	MET	3.3
36	1	1564	U	3.3
58	N2	95	PHE	3.3
16	c4	25	ASP	3.3
58	N2	9	GLN	3.3
71	o5	111	PHE	3.3
1	2	231	U	3.3
34	SR	27	ALA	3.3
35	sM	57	ASN	3.3
48	M1	135	GLY	3.3
31	D9	25	SER	3.3
16	c4	41	ARG	3.3
22	D0	117	VAL	3.3
67	O1	33	VAL	3.3
42	L5	88	ILE	3.3
11	S9	39	LYS	3.3
32	E0	29	LYS	3.3
63	N7	61	LYS	3.3
6	s4	23	LEU	3.3
3	s1	25	THR	3.3
11	s9	104	PHE	3.3
13	C1	25	VAL	3.3
17	C5	49	MET	3.3
1	2	1712	A	3.3
10	S8	151	LYS	3.3
45	L8	196	ALA	3.3
36	5	1565	G	3.3
6	S4	92	LEU	3.3
16	C4	27	PHE	3.3
21	c9	117	SER	3.3
7	S5	85	ALA	3.3
22	d0	86	ILE	3.3
23	D1	10	GLU	3.3
28	d6	55	GLU	3.3
63	N7	111	LYS	3.3
26	D4	19	ALA	3.3
83	f	122	ASP	3.3
16	c4	20	TYR	3.3
12	c0	32	HIS	3.3
21	c9	30	VAL	3.3
3	s1	52	THR	3.3
36	1	1016	C	3.3

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Mol	Chain	Res	Type	RSRZ
34	sR	67	ILE	3.3
18	c6	28	LEU	3.3
21	c9	33	TYR	3.3
34	SR	222	LEU	3.3
29	d7	32	PHE	3.3
47	M0	149	VAL	3.3
3	s1	49	ASN	3.2
18	c6	51	PRO	3.2
59	N3	32	ARG	3.2
47	M0	140	THR	3.2
15	c3	40	TYR	3.2
36	1	1238	C	3.2
81	p0	25	LEU	3.2
21	c9	73	VAL	3.2
28	D6	28	LYS	3.2
34	sR	241	PHE	3.2
63	N7	118	PHE	3.2
3	S1	190	PRO	3.2
57	n1	121	ALA	3.2
28	d6	73	TYR	3.2
2	S0	203	PHE	3.2
25	D3	130	VAL	3.2
26	D4	58	PHE	3.2
58	N2	28	PHE	3.2
63	N7	95	VAL	3.2
1	2	195	G	3.2
16	c4	85	ALA	3.2
35	sM	34	LYS	3.2
3	S1	48	VAL	3.2
8	S6	84	TYR	3.2
11	s9	149	ARG	3.2
12	C0	39	ASN	3.2
21	C9	104	VAL	3.2
74	o8	45	VAL	3.2
6	s4	69	HIS	3.2
83	f	58	HIS	3.2
7	S5	86	GLN	3.2
45	L8	214	LEU	3.2
47	m0	82	ARG	3.2
35	SM	28	SER	3.2
15	c3	83	GLU	3.2
36	5	2448	G	3.2

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Mol	Chain	Res	Type	RSRZ
63	n7	21	LYS	3.2
5	s3	50	ILE	3.2
34	sR	270	LEU	3.2
45	l8	189	LEU	3.2
58	N2	34	ALA	3.2
8	S6	97	VAL	3.2
45	l8	197	VAL	3.2
58	n2	97	SER	3.2
60	N4	73	ARG	3.2
1	2	194	U	3.2
30	D8	17	GLY	3.2
34	sR	121	MET	3.2
20	c8	137	HIS	3.2
74	o8	51	LEU	3.2
2	S0	199	PRO	3.2
4	S2	41	LEU	3.2
34	sR	48	THR	3.2
83	f	12	ASP	3.2
34	SR	212	ALA	3.2
63	N7	110	ALA	3.2
2	s0	34	GLU	3.2
15	C3	54	LEU	3.2
62	N6	92	GLY	3.2
12	C0	50	THR	3.2
12	c0	55	VAL	3.2
4	S2	151	PRO	3.2
7	S5	148	ARG	3.2
58	n2	62	VAL	3.2
34	sR	30	PRO	3.2
15	C3	53	LEU	3.2
3	S1	85	LYS	3.2
16	c4	24	ASN	3.2
40	L3	47	LEU	3.2
5	S3	77	PHE	3.2
5	s3	56	GLN	3.2
9	S7	61	PHE	3.2
10	s8	109	PHE	3.2
11	s9	184	SER	3.2
16	c4	101	ALA	3.2
26	d4	24	VAL	3.2
70	O4	110	GLU	3.2
42	l5	131	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
48	M1	102	PHE	3.2
2	S0	187	ALA	3.2
6	S4	110	ALA	3.2
21	C9	80	TYR	3.2
58	n2	16	THR	3.2
21	C9	119	LYS	3.1
45	l8	122	LYS	3.1
63	N7	70	PRO	3.1
16	C4	39	ILE	3.1
16	c4	62	LEU	3.1
40	L3	328	ILE	3.1
2	S0	196	SER	3.1
5	s3	74	GLN	3.1
6	S4	95	THR	3.1
29	D7	70	LYS	3.1
34	SR	317	THR	3.1
34	sR	53	LYS	3.1
39	L2	252	THR	3.1
2	s0	64	ILE	3.1
9	s7	32	PRO	3.1
27	D5	51	LEU	3.1
34	sR	34	LEU	3.1
20	C8	124	GLY	3.1
2	s0	99	ALA	3.1
27	d5	76	ALA	3.1
40	L3	51	ALA	3.1
42	L5	5	LYS	3.1
83	f	152	GLU	3.1
2	S0	206	ASP	3.1
14	C2	41	LEU	3.1
17	C5	127	ARG	3.1
18	C6	54	LEU	3.1
18	C6	89	LEU	3.1
40	l3	47	LEU	3.1
42	L5	131	LEU	3.1
12	C0	17	GLN	3.1
14	c2	64	SER	3.1
47	M0	95	HIS	3.1
57	N1	27	LEU	3.1
22	d0	29	THR	3.1
31	d9	11	PRO	3.1
42	l5	55	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
8	S6	187	LYS	3.1
20	C8	129	TRP	3.1
5	s3	64	ARG	3.1
8	S6	138	ALA	3.1
12	C0	45	ALA	3.1
34	sR	227	ALA	3.1
11	S9	105	LEU	3.1
14	C2	28	LEU	3.1
27	D5	69	LEU	3.1
55	M9	44	LEU	3.1
3	S1	130	SER	3.1
7	S5	164	PRO	3.1
8	S6	79	LYS	3.1
2	S0	110	TYR	3.1
48	M1	12	LEU	3.1
34	sR	263	PHE	3.1
8	S6	179	VAL	3.1
22	D0	116	VAL	3.1
34	sR	50	ASP	3.1
81	p0	16	ARG	3.1
1	6	656	G	3.1
9	S7	93	LEU	3.1
11	s9	110	GLN	3.1
26	D4	74	LEU	3.1
2	s0	166	GLY	3.1
6	S4	90	ILE	3.1
5	S3	65	ARG	3.1
22	D0	87	HIS	3.1
28	D6	85	ARG	3.1
51	M5	60	VAL	3.1
60	N4	95	SER	3.1
60	n4	95	SER	3.1
20	c8	32	LEU	3.1
63	n7	22	LYS	3.1
81	p0	18	TYR	3.1
2	S0	197	ILE	3.1
36	1	1268	G	3.1
42	L5	127	GLY	3.1
47	m0	124	GLY	3.1
5	s3	48	VAL	3.1
58	n2	66	VAL	3.1
9	S7	155	ASP	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
36	1	2205	U	3.1
22	d0	88	LYS	3.1
26	D4	68	LYS	3.1
29	D7	33	LEU	3.1
68	o2	93	ALA	3.1
42	L5	99	TYR	3.1
83	f	49	THR	3.1
35	sM	25	ILE	3.1
42	L5	247	ILE	3.1
63	n7	4	PHE	3.1
5	s3	41	VAL	3.1
11	s9	183	ALA	3.1
18	C6	5	PRO	3.1
25	D3	42	PRO	3.1
27	d5	104	ALA	3.1
4	S2	224	PHE	3.1
5	s3	76	ARG	3.1
42	L5	63	GLN	3.1
1	2	1766	A	3.1
42	L5	126	GLU	3.1
8	S6	96	SER	3.1
26	d4	72	PHE	3.1
28	d6	45	VAL	3.1
60	n4	70	LYS	3.1
83	f	82	VAL	3.1
20	C8	3	LEU	3.1
47	M0	153	ARG	3.1
13	C1	16	GLN	3.0
15	C3	139	TRP	3.0
18	c6	143	ARG	3.0
76	Q0	121	LEU	3.0
3	S1	200	ALA	3.0
6	s4	165	ALA	3.0
22	D0	119	ALA	3.0
12	C0	4	PRO	3.0
13	C1	137	PHE	3.0
36	1	1762	C	3.0
5	s3	103	GLU	3.0
9	s7	66	SER	3.0
1	2	280	U	3.0
36	1	3275	U	3.0
42	L5	159	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
81	p0	80	VAL	3.0
10	s8	66	SER	3.0
21	c9	71	VAL	3.0
1	6	231	U	3.0
36	5	2508	U	3.0
25	d3	86	PHE	3.0
33	e1	85	TYR	3.0
34	sR	27	ALA	3.0
57	n1	24	ALA	3.0
74	o8	11	PHE	3.0
8	s6	203	GLU	3.0
11	S9	163	PRO	3.0
10	S8	96	LEU	3.0
8	S6	180	THR	3.0
42	L5	55	PHE	3.0
60	n4	75	THR	3.0
11	S9	35	GLY	3.0
18	C6	142	TYR	3.0
68	o2	2	ALA	3.0
21	C9	114	VAL	3.0
5	s3	7	LYS	3.0
6	s4	134	LYS	3.0
11	S9	150	LEU	3.0
18	C6	117	LEU	3.0
19	C7	100	LEU	3.0
10	S8	95	THR	3.0
16	c4	91	THR	3.0
19	C7	126	ALA	3.0
21	C9	50	ALA	3.0
40	L3	48	GLY	3.0
7	S5	67	PRO	3.0
8	s6	214	LYS	3.0
11	S9	5	PRO	3.0
12	C0	83	PRO	3.0
3	S1	216	LYS	3.0
18	C6	118	ILE	3.0
42	L5	64	ILE	3.0
7	S5	24	VAL	3.0
5	s3	21	LEU	3.0
12	C0	32	HIS	3.0
19	c7	110	VAL	3.0
36	1	1269	U	3.0

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Mol	Chain	Res	Type	RSRZ
1	6	676	G	3.0
3	s1	90	GLU	3.0
5	s3	20	GLU	3.0
12	C0	31	LYS	3.0
21	c9	19	ALA	3.0
22	d0	82	TYR	3.0
49	M3	95	ILE	3.0
34	SR	58	VAL	3.0
34	sR	300	THR	3.0
11	S9	118	LEU	3.0
34	SR	252	LEU	3.0
8	S6	177	ARG	3.0
60	n4	73	ARG	3.0
10	S8	200	LYS	3.0
1	6	655	G	3.0
12	c0	33	GLU	3.0
12	c0	36	ASP	3.0
60	n4	130	SER	3.0
22	D0	85	ARG	3.0
3	s1	223	PHE	3.0
16	C4	120	PRO	3.0
6	S4	27	TYR	3.0
7	S5	68	ILE	3.0
60	n4	131	ALA	3.0
66	O0	10	ILE	3.0
2	s0	177	LEU	3.0
22	D0	67	THR	3.0
30	D8	5	THR	3.0
58	N2	80	THR	3.0
31	d9	34	TYR	3.0
32	E0	54	ARG	3.0
53	M7	178	ALA	3.0
54	M8	93	ILE	3.0
3	S1	184	LEU	3.0
12	C0	46	LEU	3.0
22	D0	21	LYS	3.0
71	o5	115	LYS	3.0
74	O8	57	ASN	3.0
2	S0	23	HIS	3.0
3	s1	157	GLN	3.0
9	s7	42	GLN	3.0
13	C1	78	THR	3.0

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Mol	Chain	Res	Type	RSRZ
4	S2	48	GLY	2.9
14	c2	126	TRP	2.9
22	D0	65	ILE	2.9
31	d9	19	ARG	2.9
11	S9	95	TYR	2.9
65	N9	56	ALA	2.9
12	c0	49	LEU	2.9
28	d6	19	LYS	2.9
34	SR	202	LEU	2.9
35	SM	27	LYS	2.9
6	s4	86	PHE	2.9
35	SM	50	ASN	2.9
40	L3	331	ASN	2.9
6	s4	81	THR	2.9
9	S7	74	GLN	2.9
11	S9	27	GLU	2.9
11	s9	147	MET	2.9
18	c6	11	GLY	2.9
36	1	1813	A	2.9
42	l5	38	THR	2.9
62	N6	120	GLN	2.9
4	s2	92	ALA	2.9
5	S3	75	LYS	2.9
6	s4	37	LYS	2.9
68	O2	127	ALA	2.9
9	S7	43	PHE	2.9
11	S9	185	GLY	2.9
12	c0	39	ASN	2.9
34	SR	174	ASN	2.9
48	M1	108	GLU	2.9
51	m5	147	ARG	2.9
9	S7	63	PRO	2.9
9	S7	109	VAL	2.9
12	c0	69	THR	2.9
19	C7	86	PRO	2.9
26	D4	9	THR	2.9
62	N6	116	LYS	2.9
26	d4	35	VAL	2.9
34	SR	244	ALA	2.9
55	M9	177	VAL	2.9
9	S7	92	PHE	2.9
83	f	24	SER	2.9

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Mol	Chain	Res	Type	RSRZ
28	d6	20	PRO	2.9
44	L7	26	VAL	2.9
63	n7	23	VAL	2.9
36	1	1605	A	2.9
83	f	155	ARG	2.9
21	C9	92	LYS	2.9
7	s5	153	GLY	2.9
2	s0	33	GLN	2.9
20	C8	73	MET	2.9
59	n3	3	GLY	2.9
18	C6	29	ILE	2.9
18	C6	90	VAL	2.9
34	sR	33	LEU	2.9
1	2	717	C	2.9
5	s3	31	GLU	2.9
71	O5	32	LYS	2.9
14	c2	102	GLY	2.9
1	6	506	A	2.9
66	O0	67	VAL	2.9
2	S0	161	PRO	2.9
8	S6	80	ASN	2.9
9	s7	4	PRO	2.9
11	S9	183	ALA	2.9
42	L5	119	TYR	2.9
18	C6	70	THR	2.9
6	S4	70	VAL	2.9
14	C2	136	ILE	2.9
19	C7	85	VAL	2.9
20	c8	45	LEU	2.9
45	L8	152	LEU	2.9
19	C7	116	LYS	2.9
30	d8	59	SER	2.9
31	d9	25	SER	2.9
4	S2	55	GLU	2.9
22	d0	79	TRP	2.9
42	L5	28	THR	2.9
18	c6	25	GLY	2.9
3	s1	96	LEU	2.9
17	c5	83	MET	2.9
22	d0	118	VAL	2.9
48	M1	54	VAL	2.9
79	Q3	86	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
3	s1	50	LYS	2.9
28	d6	62	TYR	2.9
33	E1	85	TYR	2.9
1	6	719	U	2.9
16	c4	57	PRO	2.9
28	d6	2	PRO	2.9
45	L8	89	GLU	2.9
16	C4	38	THR	2.9
19	C7	16	LEU	2.9
35	sM	27	LYS	2.9
42	L5	52	VAL	2.9
18	C6	64	ASP	2.9
28	D6	72	HIS	2.9
45	L8	28	HIS	2.9
20	c8	20	THR	2.9
22	d0	77	LYS	2.9
34	sR	49	GLY	2.9
55	m9	21	LYS	2.9
22	d0	65	ILE	2.9
63	n7	83	THR	2.9
81	p0	185	LEU	2.9
83	f	34	ILE	2.9
28	d6	35	ALA	2.9
45	l8	119	ALA	2.9
60	N4	68	ALA	2.9
34	SR	247	PRO	2.9
63	N7	29	HIS	2.9
6	S4	56	LEU	2.9
6	S4	153	ASN	2.9
25	d3	40	SER	2.9
55	M9	138	LEU	2.9
1	2	507	U	2.9
36	5	2505	U	2.9
63	N7	25	ILE	2.9
49	m3	111	ALA	2.9
60	N4	78	ALA	2.9
16	c4	110	LEU	2.9
29	D7	73	LEU	2.9
63	N7	40	HIS	2.9
65	n9	34	GLY	2.9
83	f	116	PRO	2.9
34	sR	92	TRP	2.8

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Mol	Chain	Res	Type	RSRZ
35	SM	37	VAL	2.8
47	M0	171	TRP	2.8
1	2	500	C	2.8
13	C1	46	LYS	2.8
2	s0	152	PRO	2.8
22	d0	63	LEU	2.8
47	M0	151	GLY	2.8
62	N6	109	LEU	2.8
21	c9	116	ILE	2.8
34	SR	188	ILE	2.8
64	n8	121	VAL	2.8
63	n7	136	PHE	2.8
25	D3	89	ASN	2.8
3	S1	106	THR	2.8
11	S9	11	THR	2.8
32	E0	41	THR	2.8
34	SR	83	ALA	2.8
36	1	1955	U	2.8
58	N2	33	TYR	2.8
8	S6	76	LEU	2.8
12	C0	19	GLY	2.8
42	L5	234	ASP	2.8
20	c8	4	VAL	2.8
22	d0	91	ILE	2.8
45	L8	232	HIS	2.8
63	n7	14	VAL	2.8
6	S4	6	LYS	2.8
34	sR	316	MET	2.8
63	N7	131	PHE	2.8
63	n7	27	LYS	2.8
5	s3	219	ALA	2.8
45	L8	114	ALA	2.8
70	O4	66	SER	2.8
25	D3	133	LEU	2.8
36	5	2538	U	2.8
40	l3	178	LEU	2.8
45	l8	238	LEU	2.8
1	2	696	C	2.8
6	s4	161	LYS	2.8
11	S9	134	ILE	2.8
26	D4	13	ILE	2.8
55	M9	183	ALA	2.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
62	N6	43	TYR	2.8
1	2	1777	G	2.8
7	S5	140	THR	2.8
11	s9	11	THR	2.8
12	C0	15	LEU	2.8
3	S1	215	VAL	2.8
20	C8	16	ARG	2.8
22	D0	97	VAL	2.8
36	5	2450	G	2.8
11	S9	31	ALA	2.8
27	d5	47	TYR	2.8
83	f	61	ALA	2.8
12	c0	29	GLN	2.8
42	L5	4	GLN	2.8
45	l8	195	SER	2.8
42	l5	41	LYS	2.8
47	M0	19	LYS	2.8
77	Q1	14	LYS	2.8
6	s4	109	PHE	2.8
7	S5	43	PHE	2.8
26	d4	57	VAL	2.8
28	d6	21	VAL	2.8
34	sR	113	VAL	2.8
1	2	232	U	2.8
21	c9	29	GLU	2.8
11	S9	159	ALA	2.8
34	SR	279	ALA	2.8
5	S3	223	LYS	2.8
45	l8	162	LEU	2.8
55	m9	180	LYS	2.8
22	d0	106	ILE	2.8
30	D8	7	VAL	2.8
34	SR	54	PHE	2.8
2	S0	80	THR	2.8
34	SR	130	THR	2.8
6	s4	36	HIS	2.8
62	N6	69	LYS	2.8
63	n7	52	LYS	2.8
72	O6	99	ARG	2.8
9	S7	126	LEU	2.8
11	s9	86	LEU	2.8
2	s0	173	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
3	S1	137	ILE	2.8
30	d8	32	PHE	2.8
36	1	2507	C	2.8
57	n1	66	ASN	2.8
10	s8	36	THR	2.8
34	sR	14	GLU	2.8
3	S1	78	ASP	2.8
45	L8	109	LEU	2.8
3	s1	68	VAL	2.8
12	c0	27	PHE	2.8
15	C3	33	VAL	2.8
28	D6	45	VAL	2.8
42	L5	53	VAL	2.8
49	m3	137	GLN	2.8
19	C7	122	ILE	2.8
49	m3	147	ILE	2.8
26	D4	20	ARG	2.8
28	D6	66	LYS	2.8
33	e1	89	LYS	2.8
34	SR	216	LYS	2.8
27	D5	65	LEU	2.8
28	d6	71	LEU	2.8
29	d7	41	LEU	2.8
22	d0	111	GLY	2.8
63	n7	91	ALA	2.8
1	2	133	U	2.8
1	2	278	U	2.8
2	S0	102	PHE	2.8
30	D8	28	VAL	2.8
34	SR	56	VAL	2.8
58	n2	54	VAL	2.8
21	c9	99	SER	2.8
83	f	47	SER	2.8
3	S1	49	ASN	2.8
8	s6	147	LEU	2.8
10	s8	95	THR	2.8
22	d0	58	LEU	2.8
21	C9	18	TYR	2.8
3	s1	65	VAL	2.8
18	c6	139	GLN	2.8
28	d6	91	ASP	2.8
40	l3	386	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
23	d1	22	ARG	2.8
5	s3	61	GLU	2.8
49	m3	136	GLU	2.8
1	2	793	A	2.8
3	s1	60	ALA	2.8
27	d5	101	TYR	2.8
45	L8	223	ALA	2.8
39	L2	224	THR	2.8
22	D0	34	LEU	2.7
1	6	1217	A	2.7
34	SR	283	LYS	2.7
36	1	1814	A	2.7
63	n7	11	ALA	2.7
65	n9	23	LYS	2.7
3	s1	20	VAL	2.7
11	S9	101	VAL	2.7
34	SR	43	ILE	2.7
51	m5	15	GLN	2.7
55	M9	72	GLU	2.7
83	f	89	GLU	2.7
6	S4	23	LEU	2.7
22	d0	26	LEU	2.7
2	s0	159	ALA	2.7
4	S2	57	PHE	2.7
14	c2	122	VAL	2.7
16	c4	23	PHE	2.7
28	D6	29	SER	2.7
30	d8	17	GLY	2.7
35	SM	42	ALA	2.7
42	L5	134	ALA	2.7
55	M9	132	PHE	2.7
59	N3	6	ALA	2.7
12	c0	43	ILE	2.7
21	c9	90	PRO	2.7
31	D9	38	ILE	2.7
36	5	2449	A	2.7
47	M0	55	ASN	2.7
63	N7	106	GLN	2.7
6	s4	245	LYS	2.7
45	L8	189	LEU	2.7
53	M7	163	LYS	2.7
55	M9	181	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
2	s0	100	GLY	2.7
8	s6	192	ALA	2.7
43	L6	135	VAL	2.7
65	n9	27	TYR	2.7
2	S0	135	GLU	2.7
4	S2	65	GLU	2.7
4	S2	178	ILE	2.7
34	sR	220	ILE	2.7
49	M3	140	SER	2.7
11	s9	74	ASN	2.7
3	S1	126	THR	2.7
8	S6	154	ARG	2.7
18	C6	47	LYS	2.7
34	sR	116	ASP	2.7
42	L5	56	THR	2.7
63	n7	115	LYS	2.7
72	o6	60	LEU	2.7
83	f	93	LEU	2.7
11	S9	122	VAL	2.7
15	C3	24	ALA	2.7
16	C4	17	ALA	2.7
16	c4	59	ALA	2.7
43	l6	3	ALA	2.7
49	M3	96	ALA	2.7
1	6	657	U	2.7
6	s4	64	ILE	2.7
16	c4	19	ILE	2.7
17	C5	51	SER	2.7
58	n2	41	ILE	2.7
18	C6	12	LYS	2.7
2	S0	72	ASP	2.7
8	S6	71	THR	2.7
34	SR	144	LEU	2.7
10	S8	67	TRP	2.7
5	s3	17	PHE	2.7
5	s3	24	PHE	2.7
2	s0	50	VAL	2.7
11	S9	98	ALA	2.7
45	l8	154	ALA	2.7
8	s6	194	LYS	2.7
21	c9	94	ILE	2.7
28	d6	30	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
31	d9	12	ARG	2.7
79	q3	3	LYS	2.7
5	s3	72	LEU	2.7
11	S9	116	LEU	2.7
36	5	3155	U	2.7
53	M7	179	GLN	2.7
22	D0	72	ASN	2.7
63	n7	122	HIS	2.7
5	S3	25	PHE	2.7
16	c4	109	GLY	2.7
34	SR	280	GLY	2.7
34	sR	186	PHE	2.7
5	s3	148	LYS	2.7
12	c0	21	VAL	2.7
2	s0	122	ILE	2.7
20	c8	145	ARG	2.7
28	D6	81	ALA	2.7
42	L5	158	ARG	2.7
3	S1	54	LEU	2.7
10	s8	58	LEU	2.7
23	D1	69	LEU	2.7
32	E0	38	LEU	2.7
58	n2	105	LEU	2.7
1	6	239	C	2.7
34	sR	54	PHE	2.7
74	o8	32	ASN	2.7
42	L5	100	ALA	2.7
74	o8	55	VAL	2.7
4	S2	69	ILE	2.7
20	c8	98	TYR	2.7
34	sR	123	ILE	2.7
49	m3	79	GLU	2.7
7	s5	198	LEU	2.7
11	s9	109	LEU	2.7
1	2	140	A	2.7
16	C4	13	VAL	2.7
28	d6	51	ARG	2.7
42	L5	135	VAL	2.7
2	S0	175	TYR	2.7
3	S1	164	ILE	2.7
3	s1	98	THR	2.7
6	s4	84	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
11	s9	8	TYR	2.7
42	L5	150	LEU	2.7
72	o6	57	LEU	2.7
8	s6	79	LYS	2.7
10	s8	24	LYS	2.7
16	C4	107	ARG	2.7
20	C8	8	GLN	2.7
21	c9	106	GLN	2.7
32	E0	6	GLY	2.7
42	L5	200	PHE	2.7
12	c0	48	SER	2.7
35	SM	41	SER	2.7
45	L8	121	SER	2.7
47	m0	96	VAL	2.7
47	m0	221	ALA	2.7
49	m3	118	GLU	2.7
81	p0	50	VAL	2.7
83	f	15	SER	2.7
12	C0	54	TYR	2.7
15	C3	66	ILE	2.7
21	C9	107	ALA	2.7
34	SR	310	ILE	2.7
35	sM	29	ASN	2.7
73	O7	88	ALA	2.7
3	s1	97	LEU	2.7
14	c2	41	LEU	2.7
25	D3	34	LEU	2.7
34	sR	23	LEU	2.7
8	S6	145	PHE	2.7
29	d7	47	PHE	2.7
3	s1	66	VAL	2.7
10	s8	102	VAL	2.7
12	C0	34	GLU	2.7
21	C9	29	GLU	2.7
22	d0	117	VAL	2.7
28	D6	65	PRO	2.7
43	l6	2	SER	2.7
47	m0	45	GLU	2.7
6	s4	169	ILE	2.7
20	C8	69	ILE	2.7
47	m0	34	TYR	2.7
58	n2	33	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
72	O6	64	SER	2.7
1	2	261	U	2.7
6	S4	159	THR	2.7
6	s4	73	ASP	2.7
64	n8	120	ASN	2.7
6	s4	39	ARG	2.7
39	L2	149	ARG	2.7
5	s3	105	MET	2.7
9	s7	5	GLN	2.7
11	s9	81	VAL	2.7
27	D5	73	GLY	2.7
2	S0	141	ILE	2.7
4	S2	64	LYS	2.6
5	s3	78	LYS	2.6
6	s4	71	LYS	2.6
16	c4	21	ALA	2.7
34	SR	318	ALA	2.7
39	l2	60	LYS	2.6
9	S7	99	LEU	2.6
53	M7	170	SER	2.6
21	c9	123	ARG	2.6
9	s7	92	PHE	2.6
16	C4	28	VAL	2.6
62	N6	90	VAL	2.6
5	s3	10	LYS	2.6
13	c1	49	ILE	2.6
21	c9	119	LYS	2.6
34	sR	257	ALA	2.6
35	sM	35	ALA	2.6
32	E0	49	LEU	2.6
48	M1	19	LEU	2.6
81	p0	64	ARG	2.6
13	c1	115	PHE	2.6
55	M9	182	ASP	2.6
1	2	711	U	2.6
5	s3	16	VAL	2.6
21	C9	72	GLY	2.6
11	S9	37	LYS	2.6
31	d9	33	LYS	2.6
34	SR	6	VAL	2.6
40	L3	79	VAL	2.6
77	q1	25	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
39	l2	72	ARG	2.6
16	c4	102	LEU	2.6
26	D4	4	ALA	2.6
34	sR	7	LEU	2.6
42	L5	79	TYR	2.6
2	s0	151	SER	2.6
34	SR	3	SER	2.6
21	C9	103	LYS	2.6
28	d6	9	GLY	2.6
33	e1	120	GLU	2.6
45	L8	92	LYS	2.6
3	S1	213	ARG	2.6
2	S0	19	ALA	2.6
16	c4	60	ALA	2.6
21	c9	65	ILE	2.6
29	D7	50	ALA	2.6
48	m1	45	PRO	2.6
57	n1	30	TYR	2.6
58	N2	108	TYR	2.6
74	o8	52	TYR	2.6
31	d9	55	PHE	2.6
45	l8	28	HIS	2.6
6	s4	75	LYS	2.6
16	c4	22	SER	2.6
16	c4	108	SER	2.6
18	C6	13	LYS	2.6
66	O0	40	LYS	2.6
18	c6	8	GLN	2.6
53	M7	172	GLN	2.6
3	s1	110	LEU	2.6
10	s8	48	THR	2.6
5	S3	87	TYR	2.6
7	s5	154	ALA	2.6
11	S9	56	ALA	2.6
11	S9	140	ILE	2.6
34	sR	292	LEU	2.6
35	SM	38	PRO	2.6
36	1	1765	U	2.6
61	N5	128	ALA	2.6
9	S7	32	PRO	2.6
9	S7	157	LYS	2.6
72	O6	100	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
2	S0	74	VAL	2.6
2	s0	181	VAL	2.6
3	S1	226	GLY	2.6
27	D5	40	VAL	2.6
33	E1	147	VAL	2.6
33	e1	80	ARG	2.6
60	n4	135	SER	2.6
62	N6	95	VAL	2.6
36	1	2567	C	2.6
63	N7	123	GLN	2.6
16	C4	76	ILE	2.6
47	M0	166	ILE	2.6
3	s1	94	LYS	2.6
23	d1	37	ALA	2.6
42	L5	30	TYR	2.6
42	L5	254	LYS	2.6
70	O4	34	HIS	2.6
42	L5	245	GLU	2.6
81	p0	188	VAL	2.6
1	6	898	A	2.6
8	S6	182	GLN	2.6
67	o1	76	SER	2.6
69	O3	56	SER	2.6
7	S5	61	TYR	2.6
10	s8	137	LYS	2.6
12	C0	14	TYR	2.6
21	c9	110	LYS	2.6
33	e1	96	LYS	2.6
49	m3	124	ILE	2.6
11	S9	6	ARG	2.6
14	C2	33	ARG	2.6
1	6	277	U	2.6
6	s4	74	GLY	2.6
19	C7	110	VAL	2.6
47	M0	50	VAL	2.6
49	m3	134	GLU	2.6
83	f	60	VAL	2.6
3	S1	186	SER	2.6
9	s7	38	LEU	2.6
19	C7	109	LEU	2.6
58	N2	13	LYS	2.6
61	N5	135	ILE	2.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
67	O1	51	LEU	2.6
76	Q0	83	LYS	2.6
3	S1	31	ASP	2.6
12	c0	41	TYR	2.6
21	C9	123	ARG	2.6
58	n2	104	ARG	2.6
61	N5	122	ALA	2.6
66	O0	59	TYR	2.6
1	2	740	A	2.6
26	D4	67	GLY	2.6
67	O1	83	GLU	2.6
11	s9	92	LYS	2.6
34	sR	261	LYS	2.6
12	C0	49	LEU	2.6
4	S2	140	ARG	2.6
18	C6	123	ARG	2.6
12	C0	63	TYR	2.6
31	D9	52	PHE	2.6
34	sR	171	SER	2.6
34	sR	254	ALA	2.6
83	f	74	SER	2.6
5	s3	220	PRO	2.6
11	s9	2	PRO	2.6
20	c8	141	THR	2.6
21	C9	37	VAL	2.6
33	e1	87	THR	2.6
34	SR	208	GLY	2.6
1	2	713	A	2.6
41	l4	186	LYS	2.6
6	S4	59	ARG	2.6
8	S6	196	ARG	2.6
17	C5	89	MET	2.6
60	N4	71	ARG	2.6
1	2	710	U	2.6
26	D4	107	GLN	2.6
38	8	81	U	2.6
39	l2	59	ALA	2.6
55	M9	131	ALA	2.6
81	p0	197	PHE	2.6
6	s4	41	SER	2.6
21	c9	125	SER	2.6
2	s0	171	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
7	S5	162	VAL	2.6
12	c0	44	LYS	2.6
35	sM	83	LYS	2.6
45	l8	118	GLU	2.6
48	m1	62	ASN	2.6
59	N3	137	VAL	2.6
63	N7	13	VAL	2.6
70	o4	101	VAL	2.6
5	S3	54	ARG	2.6
2	s0	23	HIS	2.6
20	c8	55	HIS	2.6
24	d2	126	LEU	2.6
63	n7	65	ARG	2.6
67	O1	106	THR	2.6
74	o8	73	LEU	2.6
21	c9	107	ALA	2.6
27	d5	90	LYS	2.6
45	l8	177	TYR	2.6
21	c9	112	GLY	2.6
26	D4	56	SER	2.6
3	s1	228	LEU	2.6
8	S6	197	ASN	2.6
34	sR	225	LEU	2.6
62	n6	57	LEU	2.6
47	m0	33	ILE	2.6
47	m0	86	HIS	2.6
79	q3	5	THR	2.6
3	S1	182	ALA	2.5
15	c3	25	TRP	2.5
28	d6	68	TYR	2.5
33	e1	90	LYS	2.5
3	S1	196	GLU	2.5
1	2	269	G	2.5
1	2	1685	G	2.5
25	D3	51	GLY	2.5
34	sR	124	SER	2.5
40	L3	277	SER	2.5
63	n7	26	VAL	2.5
70	o4	55	SER	2.5
4	S2	240	LEU	2.5
63	N7	81	LEU	2.5
42	L5	84	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
47	M0	159	PHE	2.5
62	n6	45	ILE	2.5
20	C8	55	HIS	2.5
34	sR	203	THR	2.5
16	C4	20	TYR	2.5
45	L8	123	GLN	2.5
83	f	145	GLU	2.5
17	c5	113	GLY	2.5
21	C9	5	SER	2.5
34	SR	213	SER	2.5
5	s3	5	ILE	2.5
9	S7	113	PRO	2.5
20	C8	125	ILE	2.5
21	c9	135	ILE	2.5
20	C8	44	ASN	2.5
36	1	1242	G	2.5
2	S0	83	GLN	2.5
3	s1	92	GLN	2.5
16	c4	68	ALA	2.5
35	SM	58	GLU	2.5
12	C0	53	GLY	2.5
8	s6	131	LYS	2.5
12	C0	43	ILE	2.5
34	sR	35	SER	2.5
74	O8	56	ILE	2.5
13	C1	136	ARG	2.5
2	s0	49	ASN	2.5
34	sR	200	ASN	2.5
36	5	1025	A	2.5
13	C1	38	ALA	2.5
14	c2	42	ALA	2.5
21	c9	17	ALA	2.5
1	6	235	G	2.5
17	C5	93	VAL	2.5
21	C9	4	VAL	2.5
45	l8	157	VAL	2.5
62	N6	85	VAL	2.5
20	c8	54	LEU	2.5
61	N5	126	LEU	2.5
5	s3	94	ARG	2.5
6	s4	24	SER	2.5
20	C8	10	SER	2.5

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Mol	Chain	Res	Type	RSRZ
3	s1	233	GLY	2.5
5	s3	137	VAL	2.5
11	S9	125	ALA	2.5
17	c5	93	VAL	2.5
27	d5	40	VAL	2.5
30	D8	43	ASN	2.5
31	D9	20	GLN	2.5
34	SR	31	ASN	2.5
45	l8	130	TYR	2.5
51	m5	6	TYR	2.5
63	N7	75	VAL	2.5
63	n7	10	VAL	2.5
76	q0	90	ASN	2.5
18	C6	17	THR	2.5
21	c9	105	LEU	2.5
19	c7	60	ARG	2.5
34	SR	302	PHE	2.5
45	L8	218	ILE	2.5
34	sR	273	ASP	2.5
53	M7	16	SER	2.5
2	s0	186	GLY	2.5
5	S3	73	VAL	2.5
12	c0	54	TYR	2.5
16	C4	79	VAL	2.5
16	C4	99	GLN	2.5
20	c8	146	ALA	2.5
21	C9	33	TYR	2.5
21	C9	55	TYR	2.5
27	D5	60	VAL	2.5
29	d7	54	VAL	2.5
45	L8	198	ALA	2.5
49	M3	33	VAL	2.5
76	Q0	85	LEU	2.5
3	s1	46	THR	2.5
5	S3	76	ARG	2.5
28	d6	67	THR	2.5
1	6	1707	A	2.5
9	S7	49	ILE	2.5
15	C3	50	ILE	2.5
34	SR	192	PHE	2.5
48	M1	131	MET	2.5
8	s6	187	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
19	C7	18	GLU	2.5
19	C7	72	LYS	2.5
34	sR	5	GLU	2.5
35	sM	49	LYS	2.5
6	S4	181	VAL	2.5
12	c0	19	GLY	2.5
45	l8	115	ALA	2.5
59	N3	3	GLY	2.5
3	S1	97	LEU	2.5
12	C0	68	LEU	2.5
31	d9	41	GLN	2.5
34	SR	183	LEU	2.5
43	L6	134	ARG	2.5
17	C5	82	ASN	2.5
23	d1	83	TRP	2.5
31	d9	5	ASN	2.5
34	sR	45	TRP	2.5
3	s1	121	ILE	2.5
9	S7	91	ILE	2.5
24	d2	61	ILE	2.5
24	D2	60	LYS	2.5
10	S8	145	ALA	2.5
47	m0	128	ARG	2.5
63	N7	82	PRO	2.5
63	N7	87	LEU	2.5
83	f	26	LEU	2.5
33	E1	93	HIS	2.5
1	6	1695	G	2.5
10	S8	62	THR	2.5
47	m0	38	LYS	2.5
65	n9	22	LYS	2.5
77	Q1	7	LYS	2.5
42	L5	124	GLU	2.5
55	M9	187	GLU	2.5
2	s0	176	LEU	2.5
11	s9	141	VAL	2.5
34	sR	47	LEU	2.5
34	sR	167	VAL	2.5
52	m6	61	ALA	2.5
1	2	173	A	2.5
1	2	492	A	2.5
1	6	225	A	2.5

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Mol	Chain	Res	Type	RSRZ
11	s9	152	SER	2.5
16	c4	65	GLN	2.5
49	m3	98	ASP	2.5
18	c6	26	LYS	2.5
22	d0	66	SER	2.5
59	N3	7	GLN	2.5
27	D5	70	LYS	2.5
72	O6	27	SER	2.5
20	c8	129	TRP	2.5
46	L9	10	ILE	2.5
55	m9	179	GLU	2.5
77	Q1	17	ARG	2.5
3	S1	225	VAL	2.5
4	S2	222	TYR	2.5
9	S7	38	LEU	2.5
12	C0	55	VAL	2.5
19	C7	115	LEU	2.5
35	sM	60	ALA	2.5
39	L2	59	ALA	2.5
45	L8	55	TYR	2.5
45	L8	210	ALA	2.5
58	N2	54	VAL	2.5
62	n6	104	LEU	2.5
63	n7	24	VAL	2.5
64	N8	73	LEU	2.5
29	d7	38	PRO	2.5
34	sR	61	PHE	2.5
42	L5	145	PHE	2.5
46	L9	190	ASP	2.5
3	S1	171	ILE	2.5
13	C1	77	SER	2.5
43	L6	88	SER	2.5
73	o7	84	SER	2.5
16	C4	135	ARG	2.5
24	D2	68	ARG	2.5
2	s0	147	THR	2.5
4	s2	90	THR	2.5
18	C6	32	ASN	2.5
5	S3	218	LEU	2.5
6	s4	44	LEU	2.5
6	s4	182	TYR	2.5
16	C4	102	LEU	2.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
42	L5	93	THR	2.5
21	C9	110	LYS	2.5
58	n2	37	LEU	2.5
65	N9	54	LEU	2.5
7	s5	82	PHE	2.5
11	s9	48	GLN	2.5
5	S3	90	ARG	2.5
5	s3	49	ILE	2.5
27	D5	71	ILE	2.5
28	D6	52	ASP	2.5
34	sR	246	SER	2.5
35	SM	45	SER	2.5
36	5	1349	G	2.5
46	L9	177	ASP	2.5
58	N2	91	ASP	2.5
18	C6	93	HIS	2.5
21	c9	115	GLU	2.5
60	N4	66	GLU	2.5
1	2	489	C	2.4
17	c5	80	MET	2.4
1	2	1776	A	2.4
2	s0	25	GLY	2.4
12	C0	42	VAL	2.4
12	c0	15	LEU	2.4
26	D4	120	GLY	2.4
27	d5	46	LYS	2.4
32	e0	62	VAL	2.4
62	N6	35	LEU	2.4
66	o0	90	VAL	2.4
9	S7	173	TYR	2.4
34	sR	78	ALA	2.4
36	1	3155	U	2.4
36	5	2501	U	2.4
45	l8	199	ALA	2.4
47	M0	181	TYR	2.4
42	L5	20	PHE	2.4
7	S5	89	ILE	2.4
11	S9	126	ARG	2.4
39	l2	250	GLN	2.4
42	l5	4	GLN	2.4
74	o8	40	GLN	2.4
11	S9	121	SER	2.4

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Mol	Chain	Res	Type	RSRZ
3	S1	127	VAL	2.4
14	C2	126	TRP	2.4
20	c8	48	LYS	2.4
21	c9	9	VAL	2.4
26	D4	117	LYS	2.4
31	D9	6	VAL	2.4
34	SR	137	LYS	2.4
49	M3	172	LEU	2.4
63	n7	12	VAL	2.4
1	2	239	C	2.4
1	6	1709	C	2.4
10	S8	109	PHE	2.4
16	c4	106	ALA	2.4
29	D7	47	PHE	2.4
47	M0	68	ALA	2.4
20	C8	141	THR	2.4
35	SM	82	THR	2.4
49	m3	112	ASN	2.4
2	s0	15	GLN	2.4
8	S6	190	GLN	2.4
14	c2	71	ILE	2.4
58	N2	38	ILE	2.4
19	C7	23	LYS	2.4
49	m3	153	ASP	2.4
2	S0	146	LEU	2.4
11	s9	36	LEU	2.4
34	SR	295	SER	2.4
39	l2	29	LEU	2.4
42	l5	87	GLY	2.4
43	L6	58	LEU	2.4
2	s0	202	TYR	2.4
21	C9	57	ARG	2.4
22	d0	89	ARG	2.4
1	2	491	C	2.4
1	6	670	U	2.4
12	c0	11	ILE	2.4
24	D2	51	GLU	2.4
2	s0	161	PRO	2.4
3	S1	174	LYS	2.4
47	m0	212	GLU	2.4
36	5	1353	U	2.4
19	C7	58	MET	2.4

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Mol	Chain	Res	Type	RSRZ
5	s3	99	VAL	2.4
6	S4	17	HIS	2.4
6	s4	9	LEU	2.4
11	s9	172	VAL	2.4
2	S0	106	SER	2.4
7	s5	157	ARG	2.4
5	s3	25	PHE	2.4
10	S8	83	TYR	2.4
11	s9	164	PHE	2.4
12	C0	48	SER	2.4
30	D8	21	SER	2.4
34	SR	35	SER	2.4
34	SR	281	TYR	2.4
42	L5	220	SER	2.4
34	sR	184	ASN	2.4
42	L5	143	LYS	2.4
52	M6	42	ASN	2.4
58	N2	11	ILE	2.4
63	N7	64	LYS	2.4
13	c1	31	THR	2.4
75	o9	11	GLN	2.4
76	Q0	120	GLN	2.4
1	6	665	U	2.4
63	n7	50	PRO	2.4
2	s0	47	VAL	2.4
5	S3	48	VAL	2.4
13	C1	116	ARG	2.4
24	d2	63	VAL	2.4
4	S2	78	ASP	2.4
8	S6	54	GLY	2.4
11	s9	47	PHE	2.4
45	l8	181	LYS	2.4
47	M0	80	SER	2.4
53	M7	130	TYR	2.4
2	S0	180	GLU	2.4
7	S5	79	ASN	2.4
3	s1	134	VAL	2.4
3	s1	192	VAL	2.4
6	S4	61	VAL	2.4
6	s4	150	PRO	2.4
11	S9	160	PRO	2.4
30	D8	56	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
47	M0	69	ARG	2.4
29	D7	72	LYS	2.4
34	sR	228	LYS	2.4
42	L5	29	ASP	2.4
42	L5	164	LYS	2.4
68	o2	62	LYS	2.4
18	C6	3	ALA	2.4
1	2	1002	G	2.4
1	6	487	G	2.4
4	S2	139	ILE	2.4
49	m3	123	ILE	2.4
2	S0	113	ARG	2.4
8	s6	212	LEU	2.4
22	d0	15	GLN	2.4
40	L3	57	VAL	2.4
8	S6	64	LYS	2.4
10	S8	53	LYS	2.4
18	C6	107	LYS	2.4
58	n2	15	PHE	2.4
74	o8	26	LYS	2.4
1	6	794	U	2.4
6	s4	103	TYR	2.4
11	s9	4	ALA	2.4
14	c2	92	ALA	2.4
56	N0	2	ALA	2.4
74	o8	3	ARG	2.4
2	s0	57	LEU	2.4
60	n4	96	LEU	2.4
68	o2	4	LEU	2.4
3	S1	93	GLY	2.4
6	S4	7	LYS	2.4
13	C1	26	LYS	2.4
12	C0	28	ASN	2.4
18	C6	11	GLY	2.4
40	L3	323	MET	2.4
81	p0	27	VAL	2.4
18	C6	66	ARG	2.4
67	o1	82	GLU	2.4
34	SR	199	ILE	2.4
17	C5	105	VAL	2.4
26	d4	2	SER	2.4
28	d6	26	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
43	L6	8	LYS	2.4
47	m0	191	LYS	2.4
57	N1	106	LEU	2.4
58	n2	76	LEU	2.4
42	L5	201	GLY	2.4
46	L9	178	GLY	2.4
48	M1	83	GLY	2.4
5	S3	62	ASN	2.4
6	s4	77	ARG	2.4
9	s7	39	ARG	2.4
28	D6	25	ASN	2.4
34	SR	319	ASN	2.4
16	c4	58	TYR	2.4
33	e1	111	GLU	2.4
1	6	234	G	2.4
10	S8	183	ILE	2.4
19	c7	69	ILE	2.4
5	s3	142	LEU	2.4
47	m0	103	LEU	2.4
77	q1	13	LEU	2.4
6	S4	89	VAL	2.4
5	s3	107	PHE	2.4
7	S5	184	PHE	2.4
16	c4	44	GLY	2.4
18	C6	60	PHE	2.4
22	d0	16	GLN	2.4
27	D5	72	GLY	2.4
47	M0	136	PHE	2.4
6	s4	82	TYR	2.4
17	C5	101	ALA	2.4
24	d2	51	GLU	2.4
2	s0	92	HIS	2.4
4	S2	88	LYS	2.4
21	C9	135	ILE	2.4
21	c9	124	ILE	2.4
49	m3	93	ILE	2.4
68	o2	6	HIS	2.4
45	L8	200	LEU	2.4
45	l8	211	LEU	2.4
47	m0	87	LEU	2.4
48	M1	112	LEU	2.4
55	M9	24	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	2	1340	U	2.4
25	D3	63	GLN	2.4
28	d6	10	ARG	2.4
4	S2	46	LYS	2.4
29	D7	57	GLU	2.4
32	E0	36	LYS	2.4
32	E0	40	TYR	2.4
45	l8	113	ALA	2.4
64	N8	111	LYS	2.4
74	o8	71	PRO	2.4
15	C3	37	ILE	2.4
32	e0	51	ASN	2.4
2	s0	168	HIS	2.4
3	s1	153	HIS	2.4
45	L8	74	THR	2.4
83	f	137	THR	2.4
1	2	483	A	2.3
6	S4	72	VAL	2.3
8	S6	153	VAL	2.3
14	C2	31	VAL	2.3
21	c9	114	VAL	2.3
42	L5	125	VAL	2.3
17	C5	12	PHE	2.3
47	m0	35	ASP	2.3
3	S1	90	GLU	2.3
12	C0	44	LYS	2.3
28	D6	3	LYS	2.3
47	M0	191	LYS	2.3
3	S1	60	ALA	2.3
12	c0	63	TYR	2.3
16	C4	40	ALA	2.3
23	d1	73	ALA	2.3
45	L8	113	ALA	2.3
63	N7	77	TYR	2.3
9	S7	58	LEU	2.3
26	D4	125	LEU	2.3
48	M1	101	ASN	2.3
51	m5	22	LEU	2.3
8	s6	195	VAL	2.3
35	sM	30	THR	2.3
34	SR	304	GLY	2.3
42	l5	204	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
45	L8	84	ARG	2.3
6	S4	99	PHE	2.3
12	c0	61	TRP	2.3
16	c4	15	GLY	2.3
21	c9	121	GLY	2.3
49	M3	98	ASP	2.3
63	n7	20	GLY	2.3
17	c5	52	LYS	2.3
11	s9	153	GLU	2.3
21	C9	23	GLN	2.3
77	Q1	25	LYS	2.3
2	s0	40	ALA	2.3
32	E0	39	LEU	2.3
35	SM	47	ALA	2.3
63	n7	132	SER	2.3
8	s6	188	ARG	2.3
17	C5	10	ARG	2.3
58	N2	76	LEU	2.3
83	f	37	ARG	2.3
2	S0	47	VAL	2.3
2	S0	181	VAL	2.3
10	s8	46	VAL	2.3
1	6	1698	G	2.3
10	s8	113	PHE	2.3
22	d0	32	LYS	2.3
26	d4	23	PHE	2.3
28	d6	11	ASN	2.3
34	sR	77	GLY	2.3
57	n1	35	LYS	2.3
67	O1	34	LYS	2.3
71	O5	3	GLY	2.3
83	f	114	LYS	2.3
5	s3	93	ASP	2.3
11	s9	6	ARG	2.3
11	s9	145	SER	2.3
22	d0	119	ALA	2.3
23	d1	53	TYR	2.3
28	D6	35	ALA	2.3
31	d9	18	SER	2.3
34	SR	57	PRO	2.3
47	m0	27	PRO	2.3
55	M9	185	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
58	N2	105	LEU	2.3
6	s4	122	LYS	2.3
53	M7	174	GLY	2.3
62	n6	58	VAL	2.3
78	q2	106	PHE	2.3
14	C2	130	THR	2.3
34	sR	251	TRP	2.3
40	l3	233	TRP	2.3
46	L9	90	MET	2.3
48	M1	53	THR	2.3
58	N2	87	ASN	2.3
58	n2	61	THR	2.3
27	D5	48	ASP	2.3
16	C4	133	ARG	2.3
34	SR	90	ARG	2.3
74	o8	39	ARG	2.3
3	s1	231	LEU	2.3
9	s7	154	LEU	2.3
15	C3	16	ILE	2.3
24	D2	65	LEU	2.3
27	D5	80	LEU	2.3
6	S4	53	LYS	2.3
11	S9	154	LYS	2.3
34	SR	2	ALA	2.3
45	l8	106	LYS	2.3
67	O1	18	LYS	2.3
3	S1	38	PHE	2.3
63	N7	101	PHE	2.3
1	6	666	U	2.3
11	S9	123	HIS	2.3
56	N0	88	HIS	2.3
19	c7	106	THR	2.3
42	L5	50	ARG	2.3
46	L9	166	ARG	2.3
14	C2	42	ALA	2.3
21	c9	27	LYS	2.3
23	D1	73	ALA	2.3
25	D3	59	ILE	2.3
34	sR	73	LEU	2.3
47	m0	74	LYS	2.3
63	n7	80	LEU	2.3
1	6	1490	C	2.3

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Mol	Chain	Res	Type	RSRZ
2	S0	107	PHE	2.3
12	C0	21	VAL	2.3
34	sR	309	VAL	2.3
47	M0	147	VAL	2.3
57	N1	126	VAL	2.3
24	d2	21	GLY	2.3
36	1	1954	G	2.3
42	L5	87	GLY	2.3
42	L5	170	GLY	2.3
11	S9	149	ARG	2.3
28	d6	17	HIS	2.3
36	5	2099	A	2.3
17	c5	78	THR	2.3
4	S2	155	ALA	2.3
6	S4	82	TYR	2.3
9	s7	153	LEU	2.3
9	s7	6	ALA	2.3
21	C9	49	ASP	2.3
22	d0	104	THR	2.3
34	SR	81	LEU	2.3
47	m0	70	ILE	2.3
47	m0	77	THR	2.3
48	M1	70	THR	2.3
51	m5	39	ALA	2.3
57	n1	34	TYR	2.3
9	S7	85	PHE	2.3
28	D6	84	VAL	2.3
52	M6	7	VAL	2.3
9	s7	53	GLY	2.3
10	s8	185	GLU	2.3
31	d9	13	ARG	2.3
55	m9	170	ARG	2.3
6	s4	142	HIS	2.3
62	N6	87	LYS	2.3
7	s5	130	ILE	2.3
11	S9	45	ILE	2.3
21	C9	76	LEU	2.3
25	d3	76	LEU	2.3
36	1	2095	G	2.3
1	2	1689	A	2.3
11	S9	130	THR	2.3
12	C0	41	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
18	C6	96	TYR	2.3
29	D7	42	ASN	2.3
34	SR	4	ASN	2.3
35	sM	82	THR	2.3
62	N6	107	THR	2.3
21	c9	54	PHE	2.3
39	l2	149	ARG	2.3
74	o8	35	GLY	2.3
22	d0	83	GLU	2.3
3	s1	154	SER	2.3
59	N3	25	CYS	2.3
1	2	1686	C	2.3
3	S1	188	LEU	2.3
27	D5	83	LEU	2.3
70	o4	51	LEU	2.3
74	o8	78	LEU	2.3
7	S5	62	VAL	2.3
7	s5	133	VAL	2.3
11	s9	114	TYR	2.3
66	O0	42	ILE	2.3
13	C1	48	ALA	2.3
15	C3	26	PHE	2.3
23	D1	32	VAL	2.3
25	d3	49	ALA	2.3
27	d5	68	ARG	2.3
33	e1	84	VAL	2.3
34	SR	113	VAL	2.3
34	sR	83	ALA	2.3
1	6	1697	G	2.3
25	D3	90	ASP	2.3
74	O8	15	THR	2.3
71	O5	45	LYS	2.3
18	C6	116	LEU	2.3
34	sR	183	LEU	2.3
8	S6	101	ILE	2.3
24	d2	86	ILE	2.3
26	D4	110	GLN	2.3
34	SR	16	HIS	2.3
40	l3	109	HIS	2.3
42	L5	61	ILE	2.3
42	L5	239	ILE	2.3
6	s4	27	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
6	s4	72	VAL	2.3
7	S5	221	ALA	2.3
18	C6	10	PHE	2.3
21	c9	11	ALA	2.3
25	d3	38	PHE	2.3
61	N5	60	TYR	2.3
63	N7	49	TYR	2.3
6	s4	85	GLY	2.3
7	s5	39	GLU	2.3
55	M9	49	THR	2.3
60	n4	133	THR	2.3
68	o2	120	THR	2.3
17	c5	109	PRO	2.3
1	2	169	A	2.3
23	D1	8	LEU	2.3
30	d8	56	LEU	2.3
34	sR	222	LEU	2.3
43	L6	55	LEU	2.3
81	p0	71	PRO	2.3
8	S6	85	ARG	2.3
16	C4	103	ARG	2.3
24	D2	27	ILE	2.3
27	D5	41	ILE	2.3
31	d9	43	PHE	2.3
34	sR	205	SER	2.3
9	s7	3	ALA	2.3
12	C0	12	HIS	2.3
33	E1	83	LYS	2.3
39	L2	235	ALA	2.3
47	m0	151	GLY	2.3
81	p0	47	GLY	2.3
36	1	1764	U	2.3
6	s4	92	LEU	2.3
8	s6	173	PRO	2.3
10	S8	141	ARG	2.3
17	c5	68	PRO	2.3
21	c9	132	LEU	2.3
4	S2	218	ILE	2.3
4	s2	178	ILE	2.3
5	S3	217	ILE	2.3
10	s8	65	PHE	2.3
25	d3	50	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
41	l4	26	PHE	2.3
58	N2	71	PHE	2.3
3	s1	235	GLY	2.2
16	C4	78	ALA	2.3
18	C6	40	GLU	2.2
24	D2	59	GLY	2.2
34	SR	80	ALA	2.3
36	5	1580	A	2.3
45	l8	250	ALA	2.3
51	m5	52	GLY	2.2
12	C0	40	LEU	2.2
12	c0	3	MET	2.2
13	C1	82	ARG	2.2
21	C9	130	ARG	2.2
34	SR	96	THR	2.2
34	sR	258	THR	2.2
36	5	2509	U	2.2
45	l8	194	THR	2.2
67	O1	97	LEU	2.2
75	O9	21	ARG	2.2
2	s0	38	PHE	2.2
16	c4	112	ILE	2.2
45	l8	176	PRO	2.2
3	S1	66	VAL	2.2
6	s4	136	VAL	2.2
10	S8	160	PHE	2.2
45	L8	203	VAL	2.2
46	L9	175	PHE	2.2
6	S4	24	SER	2.2
22	d0	73	GLY	2.2
47	M0	75	TYR	2.2
34	SR	82	SER	2.2
42	l5	205	SER	2.2
53	M7	177	ALA	2.2
1	6	1800	A	2.2
9	s7	88	ARG	2.2
10	s8	196	LEU	2.2
18	c6	132	LYS	2.2
3	S1	173	THR	2.2
19	C7	82	ASP	2.2
22	D0	69	LYS	2.2
63	N7	116	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
81	p0	67	LEU	2.2
39	L2	176	ASP	2.2
1	2	841	U	2.2
5	S3	17	PHE	2.2
21	c9	36	ILE	2.2
29	D7	32	PHE	2.2
36	1	1241	U	2.2
42	l5	190	ILE	2.2
45	L8	132	VAL	2.2
61	N5	84	PHE	2.2
11	S9	12	TYR	2.2
19	c7	62	GLN	2.2
45	L8	107	GLU	2.2
47	M0	186	GLU	2.2
60	n4	128	ALA	2.2
62	N6	93	ALA	2.2
81	p0	191	TYR	2.2
2	S0	185	ARG	2.2
28	d6	72	HIS	2.2
60	n4	72	SER	2.2
6	S4	38	LEU	2.2
11	S9	30	LEU	2.2
11	s9	180	LYS	2.2
63	n7	87	LEU	2.2
3	S1	32	ILE	2.2
5	S3	50	ILE	2.2
7	s5	145	ASP	2.2
7	S5	133	VAL	2.2
11	S9	7	THR	2.2
11	S9	129	ILE	2.2
20	c8	23	ASP	2.2
11	S9	113	VAL	2.2
14	C2	82	PRO	2.2
18	C6	69	VAL	2.2
30	D8	48	VAL	2.2
34	sR	168	THR	2.2
42	l5	37	VAL	2.2
49	M3	86	THR	2.2
54	M8	88	THR	2.2
83	f	88	ASN	2.2
1	6	1488	G	2.2
6	S4	124	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
32	e0	40	TYR	2.2
67	o1	111	GLU	2.2
63	N7	60	LYS	2.2
1	6	1705	C	2.2
2	s0	54	TRP	2.2
11	s9	118	LEU	2.2
31	D9	36	LEU	2.2
2	S0	48	ILE	2.2
12	C0	59	PHE	2.2
12	C0	73	VAL	2.2
31	D9	23	VAL	2.2
47	M0	31	ILE	2.2
52	M6	3	VAL	2.2
72	o6	93	ILE	2.2
2	S0	101	ARG	2.2
21	c9	8	ASP	2.2
3	S1	102	GLY	2.2
3	S1	227	ALA	2.2
9	s7	63	PRO	2.2
4	S2	150	GLN	2.2
18	C6	128	LYS	2.2
25	D3	31	LYS	2.2
28	D6	32	LYS	2.2
35	SM	54	PRO	2.2
48	M1	73	GLY	2.2
49	m3	144	THR	2.2
57	n1	73	GLY	2.2
62	N6	88	GLU	2.2
76	Q0	108	THR	2.2
81	p0	94	THR	2.2
1	2	908	U	2.2
7	S5	72	HIS	2.2
1	6	1433	G	2.2
15	C3	25	TRP	2.2
50	M4	60	LEU	2.2
58	N2	67	SER	2.2
83	f	76	SER	2.2
2	s0	76	ILE	2.2
7	s5	156	ARG	2.2
22	d0	108	ILE	2.2
27	D5	49	ARG	2.2
47	M0	138	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
49	m3	138	VAL	2.2
58	N2	93	ILE	2.2
8	S6	86	PRO	2.2
10	S8	167	ALA	2.2
17	C5	17	TYR	2.2
22	d0	72	ASN	2.2
33	e1	110	ALA	2.2
34	SR	227	ALA	2.2
46	l9	3	TYR	2.2
3	S1	172	LEU	2.2
28	D6	77	CYS	2.2
45	L8	186	LEU	2.2
70	o4	57	LEU	2.2
36	1	252	U	2.2
42	L5	203	HIS	2.2
6	S4	147	ILE	2.2
7	S5	98	MET	2.2
8	S6	156	PHE	2.2
16	C4	118	VAL	2.2
16	c4	92	LYS	2.2
18	c6	69	VAL	2.2
20	C8	123	ARG	2.2
24	D2	34	ILE	2.2
40	L3	336	VAL	2.2
42	L5	27	LYS	2.2
45	L8	98	ARG	2.2
48	M1	67	VAL	2.2
74	O8	5	ILE	2.2
1	6	490	C	2.2
5	s3	3	ALA	2.2
8	s6	208	TYR	2.2
36	1	1017	C	2.2
36	1	1574	C	2.2
57	n1	23	GLY	2.2
9	S7	171	ALA	2.2
20	c8	121	ALA	2.2
6	S4	98	ASN	2.2
23	D1	14	PRO	2.2
42	l5	32	GLN	2.2
45	L8	199	ALA	2.2
6	s4	164	LEU	2.2
45	l8	191	ASN	2.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
20	C8	68	ARG	2.2
1	2	909	U	2.2
3	S1	29	TRP	2.2
6	S4	226	PHE	2.2
16	c4	115	ILE	2.2
25	D3	137	LYS	2.2
63	N7	136	PHE	2.2
35	sM	26	VAL	2.2
54	M8	167	SER	2.2
81	p0	190	VAL	2.2
17	C5	123	TYR	2.2
47	M0	34	TYR	2.2
55	m9	188	ASP	2.2
65	N9	2	ALA	2.2
1	2	1796	C	2.2
5	s3	109	LEU	2.2
11	s9	97	LEU	2.2
11	s9	150	LEU	2.2
63	n7	128	GLN	2.2
70	o4	26	PRO	2.2
6	S4	11	ARG	2.2
10	s8	59	ARG	2.2
71	o5	106	LYS	2.2
4	S2	103	VAL	2.2
16	c4	83	ILE	2.2
30	D8	44	VAL	2.2
42	l5	126	GLU	2.2
43	L6	96	VAL	2.2
48	M1	125	MET	2.2
83	f	21	MET	2.2
16	c4	98	GLY	2.2
63	n7	97	SER	2.2
32	E0	2	ALA	2.2
68	o2	92	TYR	2.2
3	S1	35	PRO	2.2
3	S1	104	ASP	2.2
3	s1	74	GLN	2.2
8	s6	191	ARG	2.2
31	D9	40	ARG	2.2
34	SR	165	ASP	2.2
22	d0	36	ASN	2.2
63	N7	127	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
23	D1	82	VAL	2.2
34	SR	290	VAL	2.2
34	sR	178	VAL	2.2
34	sR	221	MET	2.2
46	L9	140	VAL	2.2
57	n1	76	ILE	2.2
67	O1	64	VAL	2.2
20	c8	37	GLY	2.2
51	M5	58	GLY	2.2
1	2	504	U	2.2
1	2	1687	U	2.2
1	2	1795	U	2.2
3	S1	51	SER	2.2
13	C1	40	LEU	2.2
47	M0	155	ALA	2.2
55	m9	181	ARG	2.2
55	m9	183	ALA	2.2
63	N7	105	SER	2.2
68	O2	51	SER	2.2
1	6	673	A	2.2
34	sR	51	ASP	2.2
65	n9	24	PRO	2.2
27	D5	64	VAL	2.2
43	L6	97	ASN	2.2
47	M0	76	MET	2.2
57	n1	48	ILE	2.2
61	N5	111	ASN	2.2
11	S9	124	HIS	2.2
23	D1	57	GLY	2.2
32	e0	52	GLY	2.2
60	n4	84	GLY	2.2
2	s0	41	ARG	2.2
5	s3	29	LEU	2.2
9	S7	44	LYS	2.2
19	C7	59	LYS	2.2
21	c9	86	ARG	2.2
11	s9	181	ALA	2.2
20	c8	15	LEU	2.2
58	N2	79	LEU	2.2
1	6	710	U	2.2
4	s2	181	SER	2.2
32	e0	44	PHE	2.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
36	1	2503	G	2.2
34	SR	76	ASP	2.2
62	N6	96	PRO	2.2
5	s3	177	MET	2.2
12	c0	78	GLU	2.2
42	L5	133	GLU	2.2
42	l5	148	ILE	2.2
11	S9	174	ARG	2.2
20	c8	123	ARG	2.2
26	D4	77	ASN	2.2
34	sR	105	GLY	2.2
42	l5	68	THR	2.2
60	N4	69	LYS	2.2
62	N6	91	ASN	2.2
63	N7	79	HIS	2.2
8	S6	178	LEU	2.1
40	L3	338	LEU	2.1
49	m3	148	ALA	2.1
58	n2	103	TYR	2.1
59	N3	54	LEU	2.1
61	N5	24	LEU	2.1
2	s0	61	ALA	2.1
66	O0	71	GLN	2.1
6	S4	43	PRO	2.1
2	s0	150	ASP	2.1
12	C0	11	ILE	2.1
12	C0	26	ASP	2.1
14	C2	30	VAL	2.1
32	e0	4	VAL	2.1
34	SR	158	PRO	2.1
34	sR	156	VAL	2.1
45	L8	185	ARG	2.1
47	m0	44	ASP	2.1
58	N2	41	ILE	2.1
1	2	285	G	2.1
1	6	677	G	2.1
18	c6	130	GLY	2.1
33	e1	78	LYS	2.1
30	d8	33	LEU	2.1
36	1	1036	A	2.1
36	1	1271	A	2.1
36	5	1813	A	2.1

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Mol	Chain	Res	Type	RSRZ
47	m0	95	HIS	2.1
29	d7	5	GLN	2.1
3	s1	219	LYS	2.1
10	s8	38	ILE	2.1
1	2	912	U	2.1
1	6	711	U	2.1
6	s4	35	PRO	2.1
18	c6	124	PRO	2.1
25	D3	39	LYS	2.1
45	l8	143	ILE	2.1
45	l8	159	PRO	2.1
47	M0	27	PRO	2.1
48	m1	173	ASP	2.1
2	S0	159	ALA	2.1
2	s0	60	ALA	2.1
3	s1	117	TRP	2.1
10	S8	44	HIS	2.1
12	C0	30	ALA	2.1
24	D2	46	TYR	2.1
24	d2	130	TYR	2.1
26	d4	9	THR	2.1
26	d4	34	ASN	2.1
27	D5	101	TYR	2.1
48	M1	147	THR	2.1
71	O5	46	THR	2.1
83	f	146	ALA	2.1
1	2	451	A	2.1
9	S7	33	GLU	2.1
33	e1	113	LYS	2.1
73	o7	85	LYS	2.1
77	Q1	15	ARG	2.1
81	p0	14	LYS	2.1
25	D3	72	VAL	2.1
27	D5	54	VAL	2.1
28	D6	88	SER	2.1
34	SR	194	GLY	2.1
74	o8	37	PRO	2.1
20	C8	101	LEU	2.1
26	d4	26	ASP	2.1
36	5	2507	C	2.1
62	N6	111	LEU	2.1
2	S0	138	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
15	C3	67	THR	2.1
47	M0	139	ARG	2.1
55	M9	172	ARG	2.1
5	s3	73	VAL	2.1
13	C1	64	VAL	2.1
15	c3	66	ILE	2.1
16	c4	74	VAL	2.1
34	SR	189	GLU	2.1
61	n5	26	VAL	2.1
64	n8	124	ILE	2.1
8	S6	69	LEU	2.1
13	C1	62	GLY	2.1
19	c7	57	LEU	2.1
42	l5	236	LEU	2.1
49	M3	51	LEU	2.1
1	2	192	U	2.1
1	6	261	U	2.1
5	S3	152	PHE	2.1
5	s3	106	LYS	2.1
18	C6	45	ARG	2.1
26	D4	41	ARG	2.1
30	D8	45	LYS	2.1
49	m3	122	LYS	2.1
58	N2	36	TYR	2.1
40	L3	177	HIS	2.1
55	M9	23	TRP	2.1
7	S5	132	VAL	2.1
23	D1	39	VAL	2.1
25	D3	53	VAL	2.1
45	l8	137	ASN	2.1
30	d8	40	ILE	2.1
47	M0	59	GLN	2.1
5	s3	71	LEU	2.1
20	C8	9	GLY	2.1
6	s4	234	PRO	2.1
9	s7	77	LEU	2.1
24	d2	26	LEU	2.1
42	L5	161	GLY	2.1
29	D7	30	SER	2.1
47	m0	198	LYS	2.1
74	O8	74	LYS	2.1
11	S9	177	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
19	C7	21	TYR	2.1
22	d0	92	ASP	2.1
36	1	2206	G	2.1
74	o8	58	ASP	2.1
1	2	768	C	2.1
1	6	275	C	2.1
5	S3	188	ILE	2.1
13	C1	117	VAL	2.1
63	n7	120	GLU	2.1
23	d1	34	ILE	2.1
27	d5	105	THR	2.1
31	d9	37	ASN	2.1
34	SR	185	GLN	2.1
34	sR	319	ASN	2.1
67	O1	108	VAL	2.1
49	m3	114	GLN	2.1
53	M7	29	THR	2.1
2	S0	184	LEU	2.1
39	l2	248	GLY	2.1
42	L5	102	GLY	2.1
48	m1	140	ARG	2.1
55	m9	165	LYS	2.1
67	O1	104	LEU	2.1
22	d0	55	PRO	2.1
26	d4	30	PRO	2.1
63	n7	82	PRO	2.1
5	s3	32	GLU	2.1
5	s3	153	ALA	2.1
10	S8	106	ALA	2.1
21	c9	10	ALA	2.1
39	l2	143	GLU	2.1
43	L6	129	GLU	2.1
45	l8	254	ASP	2.1
56	n0	2	ALA	2.1
1	2	1221	A	2.1
6	s4	102	VAL	2.1
6	s4	160	VAL	2.1
60	N4	65	GLU	2.1
34	SR	294	TRP	2.1
43	l6	149	ILE	2.1
63	n7	113	VAL	2.1
66	O0	92	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
5	S3	185	LYS	2.1
6	S4	9	LEU	2.1
6	s4	242	LYS	2.1
14	C2	34	THR	2.1
14	C2	36	LEU	2.1
25	D3	44	GLY	2.1
34	SR	89	LEU	2.1
34	SR	306	THR	2.1
34	sR	139	GLN	2.1
63	N7	27	LYS	2.1
83	f	118	GLY	2.1
11	s9	95	TYR	2.1
21	c9	91	TYR	2.1
29	d7	57	GLU	2.1
34	SR	171	SER	2.1
53	M7	2	ALA	2.1
74	o8	62	ALA	2.1
6	S4	127	LYS	2.1
10	s8	169	ILE	2.1
29	D7	46	VAL	2.1
30	d8	28	VAL	2.1
11	s9	108	ARG	2.1
11	s9	171	ARG	2.1
17	C5	128	HIS	2.1
19	C7	24	LEU	2.1
20	C8	32	LEU	2.1
31	D9	44	ARG	2.1
36	1	2207	A	2.1
36	1	2540	A	2.1
55	M9	175	GLN	2.1
2	s0	203	PHE	2.1
15	C3	61	THR	2.1
1	2	449	C	2.1
4	S2	85	PRO	2.1
5	S3	216	PRO	2.1
11	s9	169	PRO	2.1
28	d6	34	LYS	2.1
34	sR	161	LYS	2.1
44	l7	23	ALA	2.1
47	M0	11	TYR	2.1
3	s1	21	VAL	2.1
5	s3	39	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
17	c5	86	VAL	2.1
19	c7	66	VAL	2.1
3	S1	131	ASP	2.1
7	S5	219	ARG	2.1
51	m5	18	VAL	2.1
74	o8	44	LYS	2.1
42	L5	65	ILE	2.1
16	c4	45	GLY	2.1
45	l8	77	GLN	2.1
5	s3	83	THR	2.1
6	s4	91	THR	2.1
16	c4	31	THR	2.1
34	SR	203	THR	2.1
36	1	1028	U	2.1
38	4	125	U	2.1
49	m3	86	THR	2.1
8	s6	115	LYS	2.1
9	S7	4	PRO	2.1
22	d0	90	TYR	2.1
26	D4	76	TYR	2.1
35	SM	39	PRO	2.1
40	L3	115	LYS	2.1
51	M5	6	TYR	2.1
1	6	1082	C	2.1
36	1	1761	C	2.1
39	L2	134	VAL	2.1
70	O4	35	VAL	2.1
12	c0	35	ILE	2.1
18	c6	29	ILE	2.1
42	L5	235	SER	2.1
46	L9	144	ILE	2.1
2	s0	17	LEU	2.1
47	M0	61	SER	2.1
25	D3	24	TRP	2.1
42	l5	200	PHE	2.1
18	C6	14	LYS	2.1
1	2	501	U	2.1
6	S4	30	ARG	2.1
23	d1	44	ARG	2.1
24	D2	18	GLU	2.1
28	D6	62	TYR	2.1
35	SM	53	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
40	L3	174	LYS	2.1
56	N0	95	ARG	2.1
59	n3	4	ASN	2.1
62	N6	98	ASN	2.1
66	O0	64	LYS	2.1
70	o4	21	LYS	2.1
2	s0	29	VAL	2.1
17	C5	94	VAL	2.1
33	e1	102	VAL	2.1
34	SR	303	ALA	2.1
47	M0	16	PRO	2.1
1	6	492	A	2.1
2	S0	76	ILE	2.1
11	s9	134	ILE	2.1
36	5	1575	A	2.1
48	M1	122	ILE	2.1
51	m5	142	ILE	2.1
6	s4	260	GLY	2.1
11	S9	32	GLY	2.1
39	L2	223	SER	2.1
40	l3	245	GLY	2.1
62	N6	99	LEU	2.1
5	s3	40	ARG	2.0
31	D9	12	ARG	2.0
36	5	170	G	2.0
60	n4	106	GLU	2.0
17	C5	76	VAL	2.0
19	C7	124	VAL	2.0
57	N1	109	VAL	2.0
61	N5	103	TYR	2.0
34	SR	41	THR	2.0
35	sM	44	PRO	2.0
40	l3	164	THR	2.0
72	O6	22	PRO	2.0
51	m5	58	GLY	2.0
5	S3	177	MET	2.0
16	C4	41	ARG	2.0
18	c6	83	GLN	2.0
23	D1	65	SER	2.0
24	D2	71	LYS	2.0
25	D3	122	PHE	2.0
33	e1	81	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
77	Q1	8	LYS	2.0
26	D4	61	ARG	2.0
34	SR	66	HIS	2.0
34	sR	311	ARG	2.0
40	L3	80	ASP	2.0
10	S8	187	GLU	2.0
3	S1	21	VAL	2.0
10	s8	34	ALA	2.0
3	S1	99	ASN	2.0
4	S2	236	PRO	2.0
8	S6	147	LEU	2.0
16	c4	43	THR	2.0
17	C5	78	THR	2.0
34	sR	188	ILE	2.0
42	l5	36	LEU	2.0
42	l5	146	LEU	2.0
67	O1	16	LEU	2.0
79	Q3	22	LEU	2.0
83	f	151	LYS	2.0
60	N4	1	MET	2.0
81	p0	48	ARG	2.0
2	S0	157	ASP	2.0
17	C5	104	GLN	2.0
31	d9	45	GLU	2.0
35	SM	78	ASP	2.0
47	m0	62	SER	2.0
47	m0	171	TRP	2.0
3	s1	225	VAL	2.0
6	s4	105	VAL	2.0
28	D6	24	VAL	2.0
34	SR	204	ALA	2.0
63	N7	23	VAL	2.0
7	s5	137	ILE	2.0
11	S9	24	LEU	2.0
59	N3	36	ILE	2.0
83	f	25	ALA	2.0
2	S0	38	PHE	2.0
17	c5	10	ARG	2.0
21	C9	90	PRO	2.0
26	D4	90	ARG	2.0
35	sM	69	ARG	2.0
58	N2	64	THR	2.0

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Mol	Chain	Res	Type	RSRZ
67	o1	45	GLY	2.0
20	c8	11	PHE	2.0
47	m0	76	MET	2.0
1	2	377	G	2.0
8	s6	182	GLN	2.0
23	D1	33	GLN	2.0
23	D1	68	SER	2.0
25	d3	145	SER	2.0
18	c6	49	TYR	2.0
18	c6	55	VAL	2.0
18	c6	98	ASP	2.0
27	D5	67	ASP	2.0
32	e0	61	SER	2.0
21	C9	30	VAL	2.0
24	d2	62	VAL	2.0
34	SR	261	LYS	2.0
42	L5	204	VAL	2.0
47	M0	67	ALA	2.0
60	N4	41	LYS	2.0
63	n7	44	ALA	2.0
71	o5	107	LYS	2.0
1	6	1702	A	2.0
10	S8	177	GLY	2.0
11	s9	117	GLY	2.0
24	D2	37	PHE	2.0
8	s6	197	ASN	2.0
30	D8	19	THR	2.0
45	L8	188	THR	2.0
5	S3	214	GLU	2.0
21	c9	109	GLU	2.0
1	2	499	U	2.0
78	Q2	102	GLN	2.0
4	s2	64	LYS	2.0
6	S4	18	TRP	2.0
9	s7	90	VAL	2.0
18	C6	74	HIS	2.0
27	d5	60	VAL	2.0
30	D8	14	LYS	2.0
42	l5	95	TRP	2.0
48	M1	128	TYR	2.0
55	m9	51	VAL	2.0
12	c0	30	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
14	c2	32	LEU	2.0
18	c6	54	LEU	2.0
19	c7	109	LEU	2.0
28	D6	64	LEU	2.0
7	S5	75	GLY	2.0
48	m1	120	ILE	2.0
74	o8	27	ILE	2.0
49	m3	75	PHE	2.0
12	C0	3	MET	2.0
43	L6	133	GLU	2.0
50	m4	92	GLU	2.0
83	f	144	GLU	2.0
3	S1	219	LYS	2.0
7	s5	203	LYS	2.0
63	N7	67	LYS	2.0
2	S0	143	VAL	2.0
6	s4	48	LEU	2.0
6	s4	235	TYR	2.0
14	c2	31	VAL	2.0
18	C6	67	VAL	2.0
19	C7	60	ARG	2.0
19	c7	35	CYS	2.0
7	S5	175	LEU	2.0
32	e0	55	ARG	2.0
48	M1	141	ARG	2.0
67	O1	79	ARG	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
84	CH	B	75	20/21	0.95	0.21	-	44,44,50,51	0
84	CH	C	74	17/21	0.90	0.21	-	47,48,55,57	0
84	CH	C	75	20/21	0.98	0.17	-	46,47,51,51	0
84	8AN	B	76	22/23	0.96	0.22	-	42,44,45,45	0
84	8AN	C	76	22/23	0.96	0.20	-	47,47,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
84	CH	B	74	20/21	0.96	0.25	-	47,50,62,62	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
87	MG	5	3405	1/1	0.96	0.74	14.49	42,42,42,42	0
87	MG	5	3404	1/1	0.95	0.47	11.33	36,36,36,36	0
87	MG	5	3401	1/1	0.95	0.37	8.09	41,41,41,41	0
86	SPS	B	3401	23/23	0.96	0.31	2.09	37,40,54,56	0
85	ZN	q3	501	1/1	1.00	0.17	0.18	69,69,69,69	0
86	SPS	1	3401	23/23	0.94	0.24	-0.37	40,43,57,59	0
85	ZN	o7	501	1/1	0.99	0.21	-0.50	51,51,51,51	0
85	ZN	D7	101	1/1	0.81	0.21	-1.08	141,141,141,141	0
85	ZN	e1	501	1/1	0.79	0.08	-1.26	164,164,164,164	0
85	ZN	E1	501	1/1	0.97	0.05	-1.29	136,136,136,136	0
85	ZN	O7	100	1/1	0.99	0.14	-1.41	52,52,52,52	0
85	ZN	Q3	501	1/1	0.98	0.11	-1.45	77,77,77,77	0
85	ZN	q0	500	1/1	0.99	0.13	-1.76	42,42,42,42	0
85	ZN	Q0	500	1/1	0.98	0.10	-1.90	57,57,57,57	0
85	ZN	D9	101	1/1	0.98	0.08	-1.93	88,88,88,88	0
85	ZN	q2	501	1/1	0.97	0.05	-2.00	77,77,77,77	0
85	ZN	d9	101	1/1	0.99	0.09	-2.06	94,94,94,94	0
85	ZN	d7	101	1/1	0.82	0.15	-2.26	131,131,131,131	0
85	ZN	d6	500	1/1	0.97	0.06	-2.31	87,87,87,87	0
85	ZN	D6	500	1/1	0.94	0.07	-2.46	105,105,105,105	0
85	ZN	Q2	501	1/1	0.91	0.05	-3.23	76,76,76,76	0
88	OHX	1	3406	7/7	0.99	0.14	-4.40	73,73,73,73	0
88	OHX	5	3406	7/7	0.98	0.09	-4.49	95,95,95,95	0
88	OHX	1	3405	7/7	0.98	0.09	-6.82	91,91,91,91	0
87	MG	1	3402	1/1	0.93	0.65	-	21,21,21,21	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
87	MG	f	1001	1/1	0.97	0.33	-	52,52,52,52	0
87	MG	B	3402	1/1	0.96	0.32	-	40,40,40,40	0
87	MG	C	101	1/1	0.94	0.31	-	49,49,49,49	0
87	MG	5	3402	1/1	0.84	0.36	-	45,45,45,45	0
87	MG	f	1002	1/1	0.93	0.08	-	66,66,66,66	0
87	MG	1	3404	1/1	0.94	0.65	-	29,29,29,29	0
87	MG	1	3403	1/1	0.89	0.38	-	45,45,45,45	0
87	MG	5	3403	1/1	0.98	0.54	-	23,23,23,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.