



Full wwPDB EM Validation Report ⓘ

Jun 23, 2026 – 05:48 PM EDT

PDB ID : 6ORL / pdb_00006orl
EMDB ID : EMD-20174
Title : RF1 pre-accommodated 70S complex at 24 ms
Authors : Fu, Z.; Indrisiunaite, G.; Kaledhonkar, S.; Shah, B.; Sun, M.; Chen, B.; Grassucci, R.A.; Ehrenberg, M.; Frank, J.
Deposited on : 2019-04-30
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

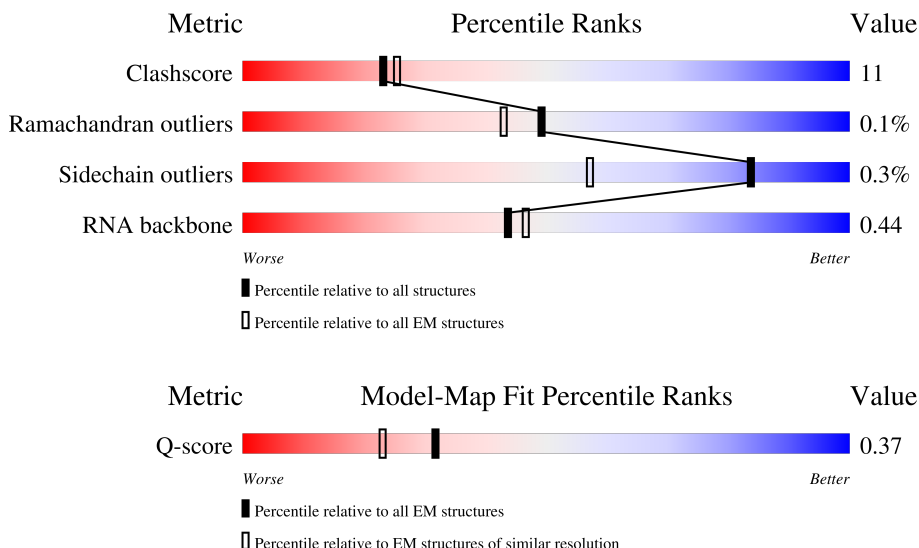
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 Å.

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



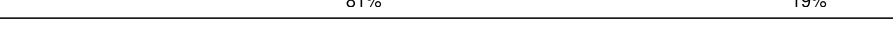
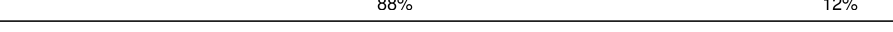
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	2903	
2	2	1534	
3	3	120	

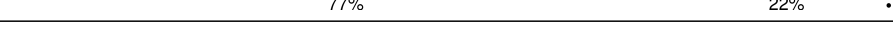
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Mol	Chain	Length	Quality of chain
4	B	271	
5	C	209	
6	D	201	
7	E	177	
8	F	175	
9	G	149	
10	J	142	
11	K	123	
12	L	144	
13	M	136	
14	N	119	
15	O	116	
16	P	114	
17	Q	117	
18	R	103	
19	S	110	
20	T	94	
21	U	103	
22	V	94	
23	W	76	
24	X	77	
25	Y	62	
26	Z	58	
27	b	56	
28	c	52	

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Mol	Chain	Length	Quality of chain
29	d	46	
30	e	64	
31	f	38	
32	g	225	
33	h	208	
34	i	205	
35	j	156	
36	k	104	
37	l	151	
38	m	129	
39	n	127	
40	o	99	
41	p	117	
42	q	123	
43	r	116	
44	s	100	
45	t	88	
46	u	82	
47	v	80	
48	w	66	
49	x	83	
50	y	86	
51	z	70	
52	5	76	
53	H	3	

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Mol	Chain	Length	Quality of chain
54	A	257	33% 67% 28%
55	4	9	67% 22% 11%

2 Entry composition

There are 57 unique types of molecules in this entry. The entry contains 146039 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	2903	Total	C	N	O	P	0	0
			62336	27816	11470	20147	2903		

- Molecule 2 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	1534	Total	C	N	O	P	0	0
			32929	14693	6041	10661	1534		

- Molecule 3 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	120	Total	C	N	O	P	0	0
			2569	1144	468	837	120		

- Molecule 4 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	B	271	Total	C	N	O	S	0	0
			2082	1288	423	364	7		

- Molecule 5 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	C	209	Total	C	N	O	S	0	0
			1565	979	288	294	4		

- Molecule 6 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	D	201	Total	C	N	O	S	0	0
			1552	974	283	290	5		

- Molecule 7 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	E	177	Total	C	N	O	S	0	0
			1410	899	249	256	6		

- Molecule 8 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	F	175	Total	C	N	O	S	0	0
			1313	826	241	244	2		

- Molecule 9 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	G	149	Total	C	N	O	S	0	0
			1111	699	197	214	1		

- Molecule 10 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	142	Total	C	N	O	S	0	0
			1129	714	212	199	4		

- Molecule 11 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	123	Total	C	N	O	S	0	0
			946	593	181	166	6		

- Molecule 12 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	144	Total	C	N	O	S	0	0
			1053	654	207	190	2		

- Molecule 13 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	136	Total	C	N	O	S	0	0
			1074	686	205	177	6		

- Molecule 14 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	119	Total	C	N	O	S	0	0
			951	588	195	163	5		

- Molecule 15 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	116	Total	C	N	O	S	0	0
			892	552	178	162			

- Molecule 16 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	114	Total	C	N	O	S	0	0
			917	574	179	163	1		

- Molecule 17 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	117	Total	C	N	O	S	0	0
			947	604	192	151			

- Molecule 18 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	R	103	Total	C	N	O	S	0	0
			816	516	153	145	2		

- Molecule 19 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	S	110	Total	C	N	O	S	0	0
			857	532	166	156	3		

- Molecule 20 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	94	Total	C	N	O	S	0	0
			746	470	140	134	2		

- Molecule 21 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms				AltConf	Trace
21	U	103	Total	C	N	O		
			788	498	148	142	0	0

- Molecule 22 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms				AltConf	Trace
22	V	94	Total	C	N	O	S	
			753	479	137	134	3	0

- Molecule 23 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms				AltConf	Trace
23	W	76	Total	C	N	O	S	
			582	360	117	104	1	0

- Molecule 24 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms				AltConf	Trace
24	X	77	Total	C	N	O	S	
			625	388	129	106	2	0

- Molecule 25 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms				AltConf	Trace
25	Y	62	Total	C	N	O	S	
			501	308	98	94	1	0

- Molecule 26 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms				AltConf	Trace
26	Z	58	Total	C	N	O	S	
			448	281	87	78	2	0

- Molecule 27 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms				AltConf	Trace
27	b	56	Total	C	N	O	S	
			444	269	94	80	1	0

- Molecule 28 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				AltConf	Trace
28	c	52	Total	C	N	O	0	0
			426	275	78	73		

- Molecule 29 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	d	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 30 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	e	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 31 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	f	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 32 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	g	225	Total	C	N	O	S	0	0
			1760	1113	316	323	8		

- Molecule 33 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	h	208	Total	C	N	O	S	0	0
			1636	1036	307	290	3		

- Molecule 34 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	i	205	Total	C	N	O	S	0	0
			1643	1026	315	298	4		

- Molecule 35 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	j	156	Total	C	N	O	S	0	0
			1152	717	217	212	6		

- Molecule 36 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	k	104	Total	C	N	O	S	0	0
			848	536	153	152	7		

- Molecule 37 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	l	151	Total	C	N	O	S	0	0
			1181	735	227	215	4		

- Molecule 38 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	m	129	Total	C	N	O	S	0	0
			979	616	173	184	6		

- Molecule 39 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	n	127	Total	C	N	O	S	0	0
			1022	634	206	179	3		

- Molecule 40 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	o	99	Total	C	N	O	S	0	0
			790	495	151	143	1		

- Molecule 41 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	p	117	Total	C	N	O	S	0	0
			877	540	174	160	3		

- Molecule 42 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	q	123	Total	C	N	O	S	0	0
			957	591	196	165	5		

- Molecule 43 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	r	116	Total	C	N	O	S	0	0
			900	558	181	158	3		

- Molecule 44 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	s	100	Total	C	N	O	S	0	0
			805	499	164	139	3		

- Molecule 45 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	t	88	Total	C	N	O	S	0	0
			714	439	144	130	1		

- Molecule 46 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	u	82	Total	C	N	O	S	0	0
			649	406	128	114	1		

- Molecule 47 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	v	80	Total	C	N	O	S	0	0
			648	411	121	113	3		

- Molecule 48 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	w	66	Total	C	N	O	S	0	0
			544	344	102	97	1		

- Molecule 49 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	x	83	Total	C	N	O	S	0	0
			663	424	126	111	2		

- Molecule 50 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	y	86	Total	C	N	O	S	0	0
			669	414	138	114	3		

- Molecule 51 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	z	70	Total	C	N	O	S	0	0
			589	366	125	97	1		

- Molecule 52 is a RNA chain called P-tRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
52	5	76	Total	C	N	O	P	S	0	0
			1627	727	296	527	76	1		

- Molecule 53 is a protein called FME-PHE-PHE.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	H	3	Total	C	N	O	S	0	0
			32	24	3	4	1		

- Molecule 54 is a protein called Peptide chain release factor 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	A	246	Total	C	N	O	S	0	0
			1929	1186	364	371	8		

- Molecule 55 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	4	9	Total	C	N	O	P	0	0
			184	83	26	66	9		

- Molecule 56 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
56	1	283	Total 283	Mg 283	0
56	2	1	Total 1	Mg 1	0
56	3	8	Total 8	Mg 8	0
56	C	1	Total 1	Mg 1	0
56	b	1	Total 1	Mg 1	0
56	i	1	Total 1	Mg 1	0

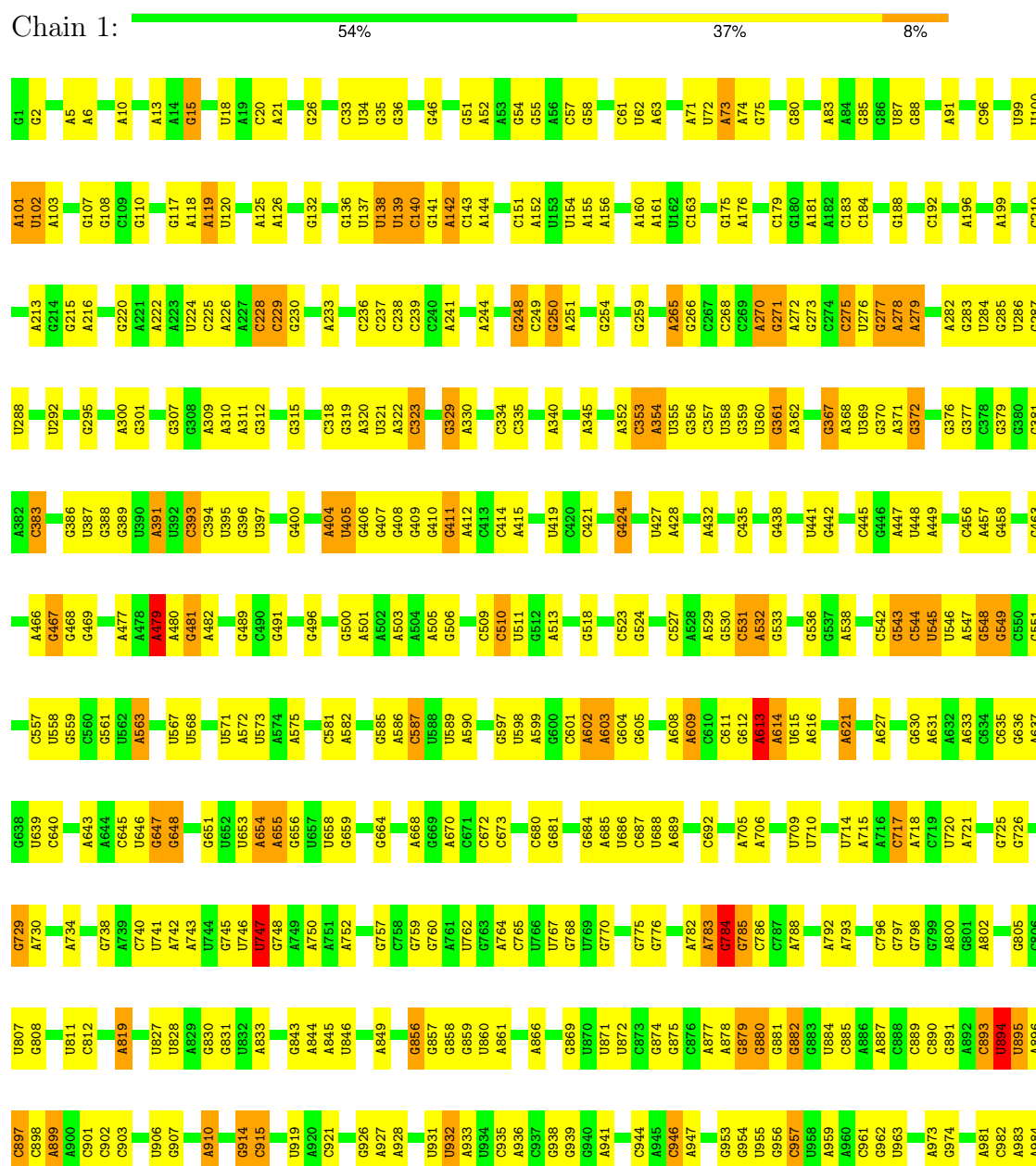
- Molecule 57 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
57	f	1	Total 1	Zn 1	0

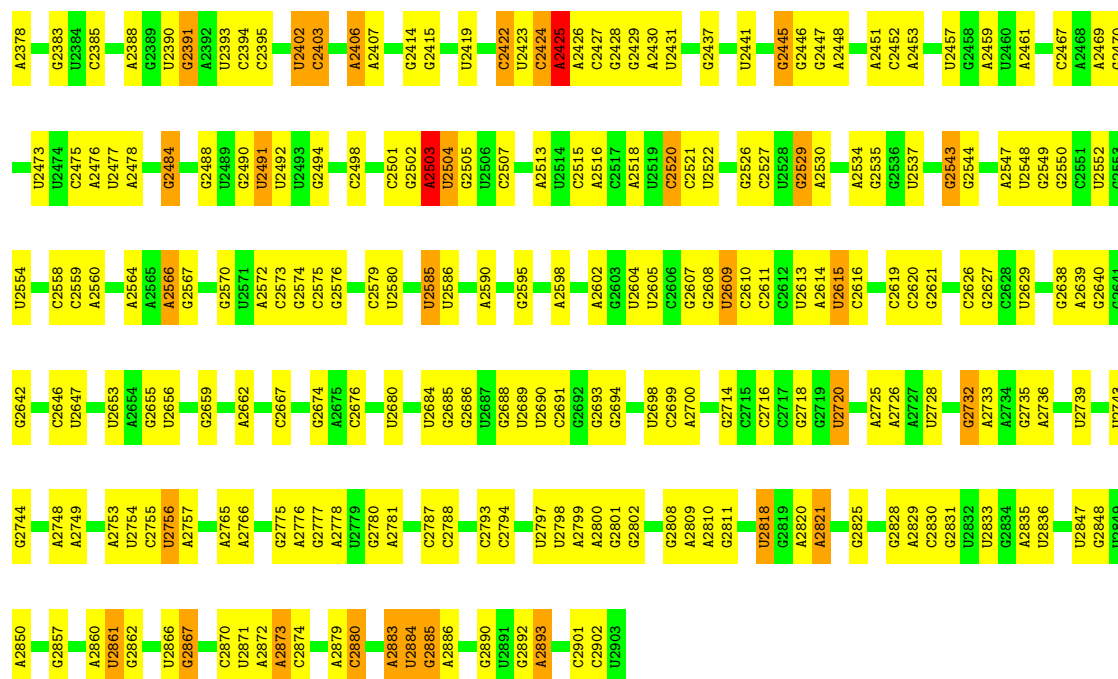
3 Residue-property plots

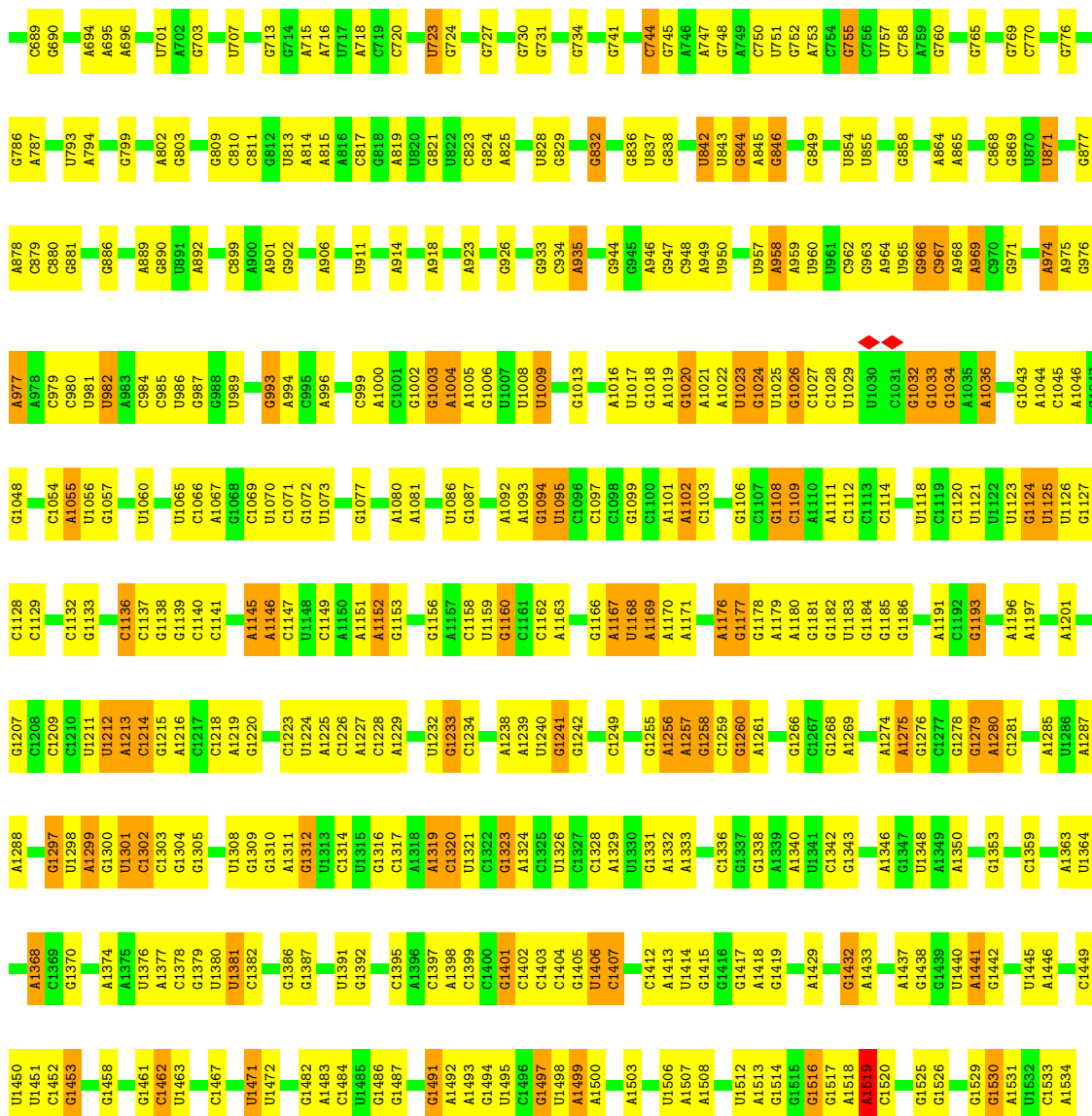
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: 23S ribosomal RNA



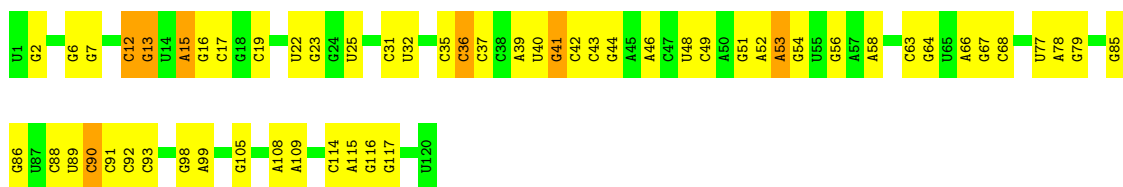
G2294	C2295	U2296	A2297	G2298	U2299	G2303	G2304	U2305	G2223	G2224	A2225	G2226	U2233	G2234	G2237	G2238	G2317	G2318	U2319	U2320	G2325	G2326	A2327	G2331	G2332	U2333	U2334	A2335	A2336	G2339	A2340	U2343	U2344	G2347	G2350	G2357	A2369	U2372	G2373	G2374	G2375	A2376	A2377							
C1934	G1935	A1936	C2036	A2037	G2038	U2039	G2040	C2043	C2047	G2048	C2055	G2056	G2057	A2058	A2059	A2060	G2061	A2062	G2063	C2064	G2069	A2070	A2071	C2072	A2077	G2087	A2088	G2093	U2099	G2100	A2101	G2109	U2110	G2111	G2112	A2114	A2115	G2116	A2117	U2118	A2119	G2120	G2121	U2122	G2123	G2125	A2126	G2127	G2128	
A1853	G1857	U1858	U1859	G1860	G1861	G1862	U1864	U1865	U1866	G1867	C1868	G1869	A1870	A1871	A1872	G1873	C1874	G1875	A1876	A1877	G1878	C1879	U1880	C1881	U1882	U1883	G1884	A1885	G1888	A1889	A1890	A1900	A1901	G1906	U1911	A1912	A1913	C1914	3TD1915	U1923	C1924	C1925	U1926	A1927	A1928	G1929	G1930			
C1934	G1935	A1936	C2036	A2037	G2038	U2039	G2040	C2043	C2047	G2048	C2055	G2056	G2057	A2058	A2059	A2060	G2061	A2062	G2063	C2064	G2069	A2070	A2071	C2072	A2077	G2087	A2088	G2093	U2099	G2100	A2101	G2110	U2111	G2112	A2114	A2115	G2116	A2117	U2118	A2119	G2120	G2121	U2122	G2123	G2125	A2126	G2127	G2128		
C2129	U2130	U2131	U2132	G2133	A2134	A2135	G2136	U2137	G2138	G2139	G2140	G2141	A2142	C2145	G2146	A2147	C2150	C2153	A2154	A2241	G2242	U2243	G2244	U2245	G2246	A2247	G2250	C2251	G2252	C2258	C2261	A2266	G2269	A2270	G2271	U2272	A2273	G2277	A2278	G2279	C2283	A2284	G2285	G2286	A2287	A2288	G2289	G2290	A2375	G2293
G2284	G2295	U2296	A2297	G2298	U2299	G2303	G2304	U2305	G2223	G2224	A2225	G2226	U2233	G2234	G2237	G2238	G2317	G2318	U2319	U2320	G2325	G2326	A2327	G2331	G2332	U2333	U2334	A2335	A2336	G2339	A2340	U2343	U2344	G2347	G2350	G2357	A2369	U2372	G2373	G2374	G2375	A2376	A2377							





• Molecule 3: 5S ribosomal RNA

Chain 3: 52% 42% 6%



• Molecule 4: 50S ribosomal protein L2

Chain B: 79% 21%





• Molecule 5: 50S ribosomal protein L3

Chain C: 84% 15%



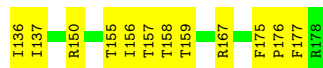
• Molecule 6: 50S ribosomal protein L4

Chain D: 81% 18%



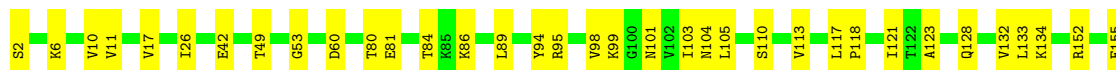
• Molecule 7: 50S ribosomal protein L5

Chain E: 74% 26%



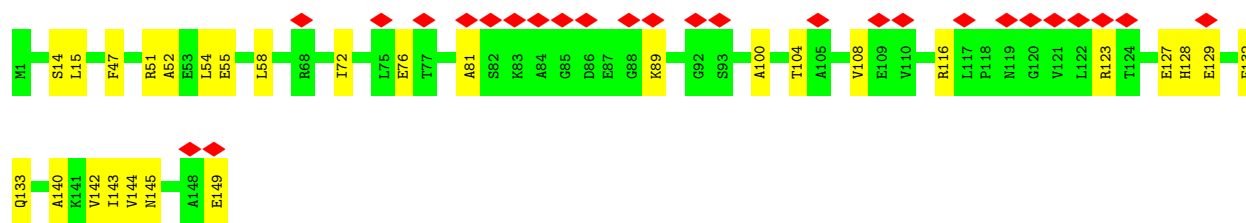
• Molecule 8: 50S ribosomal protein L6

Chain F: 77% 23%



• Molecule 9: 50S ribosomal protein L9

Chain G: 17% 81% 19%



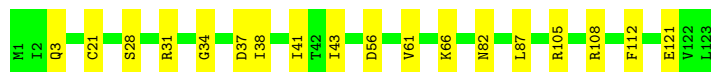
- Molecule 10: 50S ribosomal protein L13

Chain J: 89% 11%



- Molecule 11: 50S ribosomal protein L14

Chain K: 85% 15%



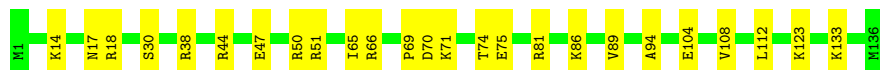
- Molecule 12: 50S ribosomal protein L15

Chain L: 85% 15%



- Molecule 13: 50S ribosomal protein L16

Chain M: 82% 18%



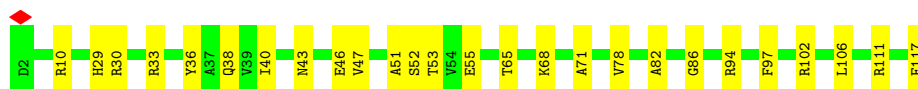
- Molecule 14: 50S ribosomal protein L17

Chain N: 87% 13%



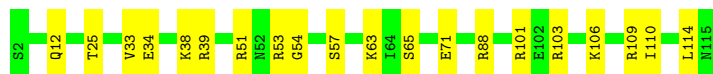
- Molecule 15: 50S ribosomal protein L18

Chain O: 78% 22%



- Molecule 16: 50S ribosomal protein L19

Chain P:  82% 18%



- Molecule 17: 50S ribosomal protein L20

Chain Q:  81% 19%



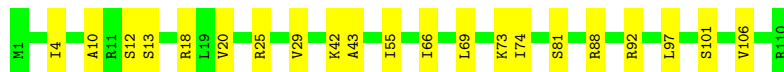
- Molecule 18: 50S ribosomal protein L21

Chain R:  82% 15% . .



- Molecule 19: 50S ribosomal protein L22

Chain S:  81% 19%



- Molecule 20: 50S ribosomal protein L23

Chain T:  88% 12%



- Molecule 21: 50S ribosomal protein L24

Chain U:  77% 22% .



- Molecule 22: 50S ribosomal protein L25

Chain V:  82% 18%



- Molecule 23: 50S ribosomal protein L27

Chain W:  84% 16%



- Molecule 24: 50S ribosomal protein L28

Chain X:  70% 30%



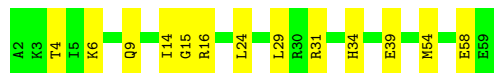
- Molecule 25: 50S ribosomal protein L29

Chain Y:  82% 18%



- Molecule 26: 50S ribosomal protein L30

Chain Z:  78% 22%



- Molecule 27: 50S ribosomal protein L32

Chain b:  73% 27%



- Molecule 28: 50S ribosomal protein L33

Chain c:  69% 31%



- Molecule 29: 50S ribosomal protein L34

Chain d:  85% 15%



- Molecule 30: 50S ribosomal protein L35

Chain e:  81% 17%



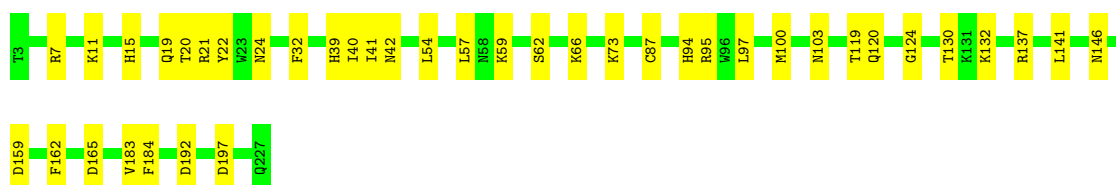
- Molecule 31: 50S ribosomal protein L36

Chain f:  71% 29%



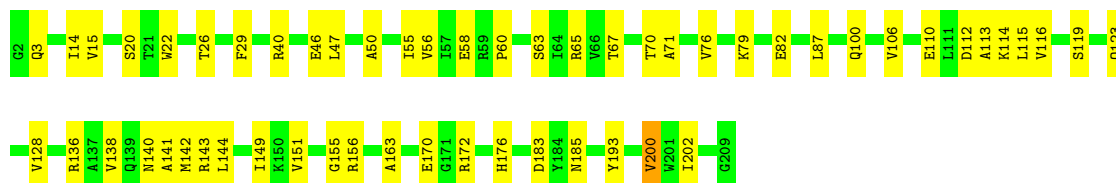
- Molecule 32: 30S ribosomal protein S2

Chain g:  82% 18%



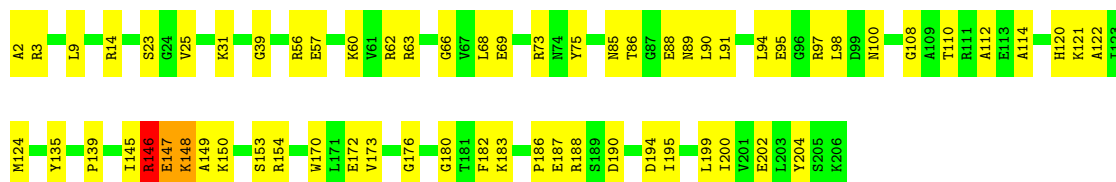
- Molecule 33: 30S ribosomal protein S3

Chain h:  74% 26%



- Molecule 34: 30S ribosomal protein S4

Chain i:  69% 30%



- Molecule 35: 30S ribosomal protein S5

Chain j:  78% 22%



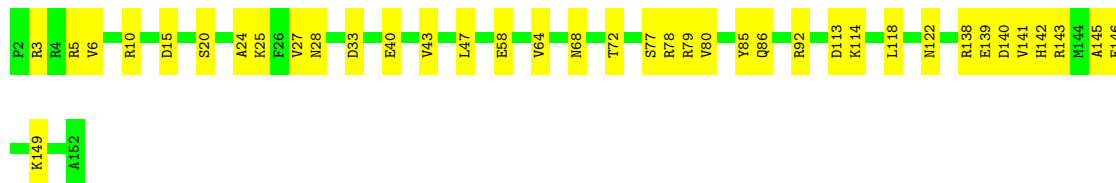
- Molecule 36: 30S ribosomal protein S6

Chain k:  75% 25%



- Molecule 37: 30S ribosomal protein S7

Chain l:  75% 25%



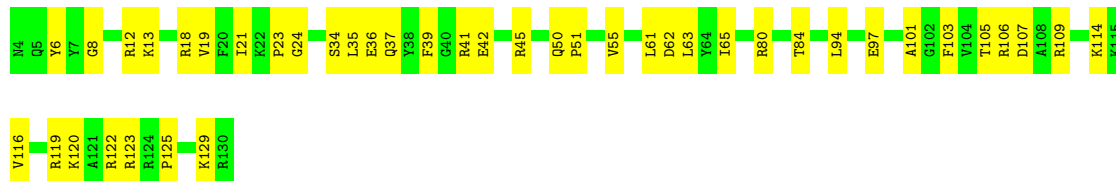
- Molecule 38: 30S ribosomal protein S8

Chain m:  80% 20%



- Molecule 39: 30S ribosomal protein S9

Chain n:  67% 33%



- Molecule 40: 30S ribosomal protein S10

Chain o:  67% 33%

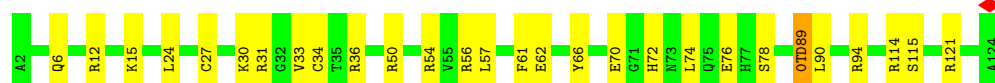
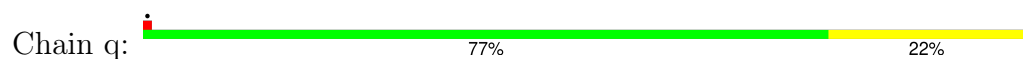


- Molecule 41: 30S ribosomal protein S11

Chain p:  76% 24%



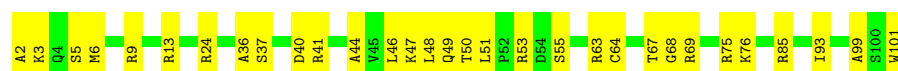
- Molecule 42: 30S ribosomal protein S12



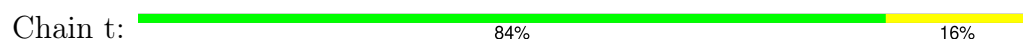
- Molecule 43: 30S ribosomal protein S13



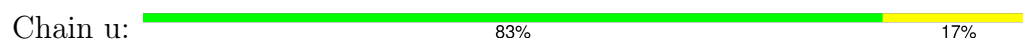
- Molecule 44: 30S ribosomal protein S14



- Molecule 45: 30S ribosomal protein S15



- Molecule 46: 30S ribosomal protein S16



- Molecule 47: 30S ribosomal protein S17



- Molecule 48: 30S ribosomal protein S18



- Molecule 49: 30S ribosomal protein S19

Chain x:  78% 22%



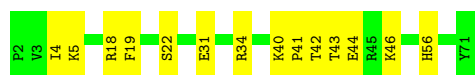
- Molecule 50: 30S ribosomal protein S20

Chain y:  84% 16%

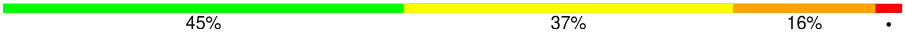


- Molecule 51: 30S ribosomal protein S21

Chain z:  80% 20%



- Molecule 52: P-tRNA

Chain 5:  45% 37% 16%



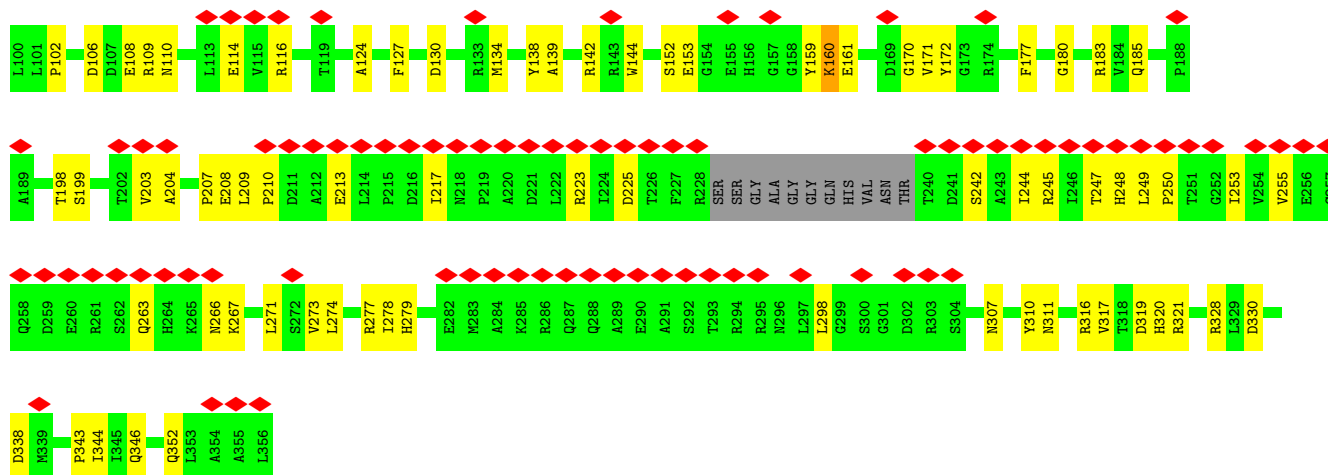
- Molecule 53: FME-PHE-PHE

Chain H:  33% 67%



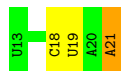
- Molecule 54: Peptide chain release factor 1

Chain A:  33% 67% 28%



- Molecule 55: mRNA

Chain 4: 



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	48706	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	41.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.147	Depositor
Minimum map value	-0.667	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.138	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	424.96, 424.96, 424.96	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.66, 1.66, 1.66	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: OMG, 0TD, PSU, OMC, UR3, ZN, 2MG, 3TD, 6MZ, FME, 4OC, MA6, 5MU, 7MG, G7M, MG, 1MG, 5MC, H2U, OMU, 4SU, 2MA

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	1	0.54	0/69286	0.44	6/108087 (0.0%)
2	2	0.37	0/36590	0.38	0/57074
3	3	0.41	0/2872	0.41	0/4478
4	B	0.51	0/2121	0.61	0/2852
5	C	0.54	0/1586	0.61	0/2134
6	D	0.49	0/1571	0.56	0/2113
7	E	0.31	0/1434	0.54	0/1926
8	F	0.32	0/1333	0.46	0/1805
9	G	0.25	0/1122	0.53	0/1515
10	J	0.55	0/1152	0.56	0/1551
11	K	0.51	0/955	0.61	0/1279
12	L	0.47	0/1062	0.57	0/1413
13	M	0.49	0/1093	0.56	0/1460
14	N	0.52	0/964	0.65	0/1289
15	O	0.34	0/902	0.53	0/1209
16	P	0.48	0/929	0.55	0/1242
17	Q	0.58	0/960	0.60	0/1278
18	R	0.56	0/829	0.70	0/1107
19	S	0.55	0/864	0.52	0/1156
20	T	0.44	0/752	0.52	0/1005
21	U	0.43	0/796	0.57	0/1062
22	V	0.38	0/766	0.50	0/1025
23	W	0.51	0/589	0.56	0/779
24	X	0.46	0/635	0.54	0/848
25	Y	0.40	0/502	0.48	0/667
26	Z	0.50	0/452	0.57	0/605
27	b	0.55	0/450	0.59	0/599
28	c	0.37	0/433	0.50	0/576
29	d	0.58	0/380	0.62	0/498
30	e	0.53	0/513	0.64	0/676
31	f	0.45	0/303	0.59	0/397

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	g	0.30	0/1791	0.55	2/2413 (0.1%)
33	h	0.30	0/1663	0.48	0/2241
34	i	0.28	0/1665	0.50	0/2227
35	j	0.35	0/1165	0.55	0/1568
36	k	0.30	0/867	0.54	0/1171
37	l	0.27	0/1195	0.50	0/1602
38	m	0.31	0/989	0.47	0/1326
39	n	0.29	0/1034	0.57	0/1375
40	o	0.27	0/800	0.50	0/1082
41	p	0.31	0/893	0.47	0/1205
42	q	0.33	0/960	0.55	0/1286
43	r	0.28	0/909	0.57	0/1215
44	s	0.27	0/817	0.45	0/1088
45	t	0.33	0/722	0.51	0/964
46	u	0.28	0/659	0.55	0/884
47	v	0.27	0/657	0.49	0/881
48	w	0.29	0/553	0.45	0/743
49	x	0.23	0/680	0.46	0/915
50	y	0.26	0/675	0.43	0/895
51	z	0.28	0/597	0.50	0/792
52	5	0.30	0/1704	0.39	0/2654
53	H	0.52	0/23	0.45	0/29
54	A	0.27	0/1959	0.59	1/2638 (0.0%)
55	4	0.29	0/203	0.46	0/312
All	All	0.46	0/157376	0.45	9/235211 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	2425	A	C4'-C3'-O3'	7.71	120.97	109.40
1	1	784	G	C4'-C3'-O3'	6.48	122.73	113.00
1	1	894	U	C4'-C3'-O3'	6.17	118.65	109.40
1	1	1584	U	C4'-C3'-O3'	5.83	118.14	109.40
1	1	613	A	C4'-C3'-O3'	5.56	117.74	109.40
54	A	160	LYS	N-CA-C	-5.42	105.27	111.07
32	g	87	CYS	CA-C-N	5.26	131.59	121.54
32	g	87	CYS	C-N-CA	5.26	131.59	121.54
1	1	479	A	C2'-C3'-O3'	-5.07	101.89	109.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	62336	0	31367	688	0
2	2	32929	0	16588	440	0
3	3	2569	0	1301	28	0
4	B	2082	0	2154	41	0
5	C	1565	0	1616	24	0
6	D	1552	0	1619	30	0
7	E	1410	0	1444	36	0
8	F	1313	0	1358	25	0
9	G	1111	0	1148	18	0
10	J	1129	0	1162	12	0
11	K	946	0	1023	12	0
12	L	1053	0	1129	18	0
13	M	1074	0	1157	18	0
14	N	951	0	994	12	0
15	O	892	0	923	19	0
16	P	917	0	962	17	0
17	Q	947	0	1019	16	0
18	R	816	0	839	19	0
19	S	857	0	922	13	0
20	T	746	0	811	8	0
21	U	788	0	843	16	0
22	V	753	0	780	10	0
23	W	582	0	599	8	0
24	X	625	0	652	16	0
25	Y	501	0	531	8	0
26	Z	448	0	488	8	0
27	b	444	0	458	12	0
28	c	426	0	464	8	0
29	d	377	0	418	7	0
30	e	504	0	572	10	0
31	f	302	0	340	11	0
32	g	1760	0	1787	25	0
33	h	1636	0	1710	38	0
34	i	1643	0	1707	50	0
35	j	1152	0	1196	20	0
36	k	848	0	846	16	0
37	l	1181	0	1238	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
38	m	979	0	1031	15	0
39	n	1022	0	1070	35	0
40	o	790	0	831	24	0
41	p	877	0	887	18	0
42	q	957	0	1017	23	0
43	r	900	0	965	23	0
44	s	805	0	844	29	0
45	t	714	0	734	11	0
46	u	649	0	666	12	0
47	v	648	0	691	17	0
48	w	544	0	560	7	0
49	x	663	0	688	15	0
50	y	669	0	719	10	0
51	z	589	0	629	10	0
52	5	1627	0	831	20	0
53	H	32	0	28	9	0
54	A	1929	0	1902	51	0
55	4	184	0	95	2	0
56	1	283	0	0	0	0
56	2	1	0	0	0	0
56	3	8	0	0	0	0
56	C	1	0	0	0	0
56	b	1	0	0	0	0
56	i	1	0	0	0	0
57	f	1	0	0	0	0
All	All	146039	0	98353	1812	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (1812) 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)
18:R:38:VAL:O	18:R:54:VAL:HG23	1.50	1.11
1:1:1380:G:H2'	1:1:1381:G:H8	1.18	1.07
1:1:1350:C:H2'	1:1:1351:C:C6	1.90	1.06
1:1:2451:A:C2	53:H:78:PHE:CE1	2.50	1.00
54:A:153:GLU:HG2	54:A:159:TYR:CD1	1.97	0.99
34:i:121:LYS:O	34:i:146:ARG:HD3	1.63	0.99
1:1:404:A:O2'	1:1:405:U:P	2.23	0.97
1:1:1380:G:C4	1:1:1381:G:N7	2.32	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:647:G:C8	1:1:647:G:H5''	2.01	0.94
1:1:1380:G:C5	1:1:1381:G:N7	2.37	0.92
1:1:1314:C:P	1:1:1332:G:H5''	2.11	0.91
1:1:893:C:O2'	1:1:894:U:O4'	1.89	0.89
1:1:1380:G:H2'	1:1:1381:G:C8	2.07	0.89
1:1:1350:C:C4	1:1:1351:C:N4	2.40	0.88
1:1:647:G:H5''	1:1:647:G:H8	1.40	0.87
54:A:153:GLU:HG2	54:A:159:TYR:HD1	1.39	0.84
1:1:404:A:O2'	1:1:405:U:OP1	1.94	0.83
2:2:677:U:H3	2:2:713:G:H22	1.29	0.81
1:1:1380:G:C6	1:1:1381:G:C6	2.68	0.81
1:1:2425:A:H5''	1:1:2427:C:O4'	1.80	0.81
1:1:99:U:HO2'	1:1:100:U:H5	1.28	0.80
54:A:153:GLU:HA	54:A:159:TYR:HA	1.64	0.80
1:1:613:A:H4'	1:1:614:A:OP1	1.81	0.78
1:1:2063:C:H1'	53:H:79:PHE:HB2	1.64	0.78
6:D:147:LEU:HD11	6:D:170:ARG:HD2	1.65	0.78
8:F:155:GLU:HG2	8:F:157:TYR:H	1.49	0.77
1:1:270:A:H5'	1:1:367:G:H22	1.49	0.77
1:1:729:G:N3	1:1:729:G:H2'	1.97	0.77
54:A:153:GLU:HG2	54:A:159:TYR:CE1	2.20	0.76
1:1:1380:G:C4	1:1:1381:G:C8	2.74	0.76
1:1:1349:C:C4	1:1:1350:C:N4	2.54	0.75
2:2:1214:C:H41	54:A:328:ARG:NH1	1.85	0.74
1:1:1350:C:H2'	1:1:1351:C:H6	1.50	0.74
1:1:783:A:H2'	1:1:783:A:N3	2.01	0.74
1:1:585:G:N7	17:Q:6:ARG:NH1	2.37	0.73
40:o:9:ARG:HB2	40:o:99:GLN:HB3	1.70	0.73
7:E:134:GLU:HG2	7:E:136:ILE:H	1.54	0.71
1:1:2451:A:C2	53:H:78:PHE:HE1	2.05	0.70
1:1:740:C:H42	1:1:757:G:H1	1.39	0.70
2:2:2:A:N1	2:2:627:G:N2	2.40	0.70
9:G:72:ILE:HD11	9:G:108:VAL:HG23	1.73	0.70
2:2:1214:C:H41	54:A:328:ARG:HH12	1.39	0.70
2:2:1255:G:O2'	2:2:1258:G:N3	2.24	0.69
2:2:531:U:OP2	55:4:21:A:N6	2.26	0.69
1:1:1568:G:N7	4:B:28:LYS:NZ	2.40	0.69
1:1:2059:A:N1	1:1:2505:G:N2	2.41	0.69
2:2:1:A:N6	2:2:627:G:N3	2.40	0.69
36:k:46:GLN:HE22	36:k:55:HIS:HB3	1.58	0.69
1:1:544:C:H2'	1:1:545:U:H4'	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:1257:A:H3'	2:2:1258:G:H8	1.57	0.68
54:A:209:LEU:HD22	54:A:279:HIS:HB2	1.76	0.68
2:2:79:G:N7	2:2:87:C:N4	2.41	0.68
2:2:1346:A:H61	2:2:1374:A:H3'	1.59	0.68
36:k:35:LYS:HB2	36:k:65:GLU:HB3	1.75	0.67
1:1:1313:U:H4'	1:1:1332:G:H4'	1.77	0.67
1:1:1380:G:O6	1:1:1381:G:O6	2.13	0.67
1:1:1913:A:N1	2:2:1491:G:N2	2.43	0.66
2:2:1081:A:N7	35:j:52:LYS:NZ	2.42	0.66
2:2:201:G:H1	2:2:216:U:H3	1.43	0.66
2:2:1120:C:H42	2:2:1153:G:H1	1.41	0.66
2:2:73:C:N3	2:2:74:A:N6	2.44	0.66
1:1:1041:G:O2'	1:1:1115:G:N2	2.28	0.66
2:2:1239:A:H62	2:2:1299:A:H62	1.42	0.66
33:h:14:ILE:HG22	33:h:15:VAL:HG13	1.77	0.66
1:1:1789:A:OP2	4:B:221:ARG:NH1	2.29	0.66
1:1:1350:C:C2	1:1:1381:G:N2	2.63	0.66
1:1:1728:C:O2'	1:1:1731:G:N2	2.28	0.66
1:1:383:C:N3	1:1:391:A:N6	2.43	0.66
1:1:2279:G:N7	23:W:14:ARG:NH2	2.43	0.65
2:2:823:C:HO2'	38:m:2:SER:N	1.94	0.65
18:R:79:ARG:HG3	18:R:80:ARG:HE	1.60	0.65
1:1:1449:G:H1	1:1:1462:C:H42	1.45	0.65
1:1:2886:A:N3	27:b:40:ARG:NH2	2.45	0.65
1:1:630:G:N2	1:1:633:A:OP2	2.28	0.65
1:1:1380:G:C6	1:1:1381:G:C5	2.84	0.65
2:2:504:C:H42	2:2:541:G:H1	1.44	0.65
2:2:1128:C:O2'	2:2:1147:C:N3	2.30	0.65
1:1:278:A:N1	1:1:361:G:O2'	2.30	0.65
1:1:468:G:N7	29:d:39:ARG:NH2	2.39	0.64
1:1:1554:U:N3	1:1:1634:A:N7	2.45	0.64
2:2:362:G:N2	2:2:365:U:OP2	2.30	0.64
1:1:871:U:H3	1:1:906:U:H3	1.44	0.64
1:1:1482:G:N2	1:1:1483:G:O6	2.31	0.64
1:1:323:C:H2'	6:D:163:ASN:ND2	2.12	0.64
1:1:1482:G:N2	1:1:1507:C:N3	2.45	0.64
2:2:1077:G:N2	2:2:1080:A:OP2	2.30	0.64
3:3:43:C:O2	7:E:92:ARG:NH2	2.30	0.64
30:e:31:HIS:HD2	30:e:32:ILE:HG12	1.61	0.64
2:2:664:G:H22	2:2:741:G:H1	1.45	0.64
2:2:949:A:H61	2:2:1232:U:H3	1.44	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2646:C:H42	1:1:2674:G:H1	1.44	0.64
31:f:25:VAL:HB	31:f:35:GLN:HB2	1.79	0.64
1:1:1776:G:H5''	1:1:1776:G:C8	2.32	0.64
2:2:933:G:O6	37:l:3:ARG:NH2	2.29	0.64
1:1:2857:G:N2	1:1:2860:A:OP2	2.31	0.64
5:C:14:ILE:HG23	16:P:12:GLN:HE22	1.63	0.64
35:j:13:GLU:HG2	35:j:39:VAL:HG12	1.80	0.64
35:j:55:GLU:HG3	35:j:57:PRO:HD2	1.79	0.64
3:3:116:G:H5'	15:O:55:GLU:HG2	1.78	0.63
47:v:76:VAL:HG23	47:v:77:ARG:HG2	1.80	0.63
54:A:153:GLU:CG	54:A:159:TYR:CD1	2.80	0.63
1:1:2395:C:N3	1:1:2422:C:N4	2.46	0.63
22:V:72:VAL:HA	22:V:94:ALA:H	1.62	0.63
1:1:228:C:N4	1:1:2407:A:N3	2.47	0.63
1:1:1405:U:H3	1:1:1597:A:H2	1.46	0.63
15:O:29:HIS:HB3	15:O:36:TYR:HB2	1.81	0.63
1:1:72:U:O4	25:Y:58:ASN:ND2	2.32	0.63
9:G:76:GLU:HA	9:G:142:VAL:HG13	1.80	0.63
2:2:254:G:N2	47:v:18:GLU:OE2	2.32	0.63
4:B:140:THR:HG22	4:B:163:GLN:HG2	1.81	0.63
54:A:153:GLU:CG	54:A:159:TYR:CE1	2.81	0.63
1:1:647:G:H3'	1:1:648:G:H8	1.64	0.63
2:2:1106:G:H5''	33:h:172:ARG:HB3	1.81	0.63
2:2:468:A:OP2	2:2:470:C:N4	2.32	0.63
37:l:27:VAL:HG12	37:l:43:VAL:HG21	1.80	0.63
1:1:819:A:OP2	1:1:1187:G:N2	2.32	0.62
1:1:13:A:O2'	1:1:15:G:N7	2.31	0.62
1:1:684:G:O2'	1:1:788:A:N7	2.32	0.62
1:1:1350:C:N3	1:1:1381:G:N2	2.34	0.62
2:2:618:C:N4	2:2:621:A:OP2	2.32	0.62
1:1:563:A:N3	17:Q:37:GLN:NE2	2.48	0.62
2:2:1167:A:O2'	2:2:1169:A:N7	2.31	0.62
27:b:30:VAL:HG22	27:b:37:LYS:HG2	1.82	0.62
2:2:501:C:OP1	42:q:114:ARG:NH2	2.32	0.62
32:g:183:VAL:HB	32:g:197:ASP:H	1.64	0.62
2:2:187:G:N2	2:2:190:A:OP2	2.32	0.62
1:1:1314:C:OP1	1:1:1332:G:H5''	1.98	0.62
2:2:4:U:H2'	2:2:5:U:H2'	1.82	0.62
1:1:879:G:OP2	1:1:893:C:N4	2.32	0.62
2:2:615:G:C5	2:2:626:G:C6	2.88	0.62
2:2:1414:U:H2'	2:2:1415:G:H8	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:i:147:GLU:O	34:i:149:ALA:N	2.33	0.61
1:1:1365:A:N7	1:1:1808:A:N6	2.47	0.61
1:1:2530:A:N7	8:F:172:LYS:NZ	2.48	0.61
4:B:144:VAL:HB	4:B:154:LEU:HB2	1.81	0.61
2:2:1073:U:O2	32:g:103:ASN:ND2	2.33	0.61
2:2:363:A:N6	42:q:27:CYS:SG	2.74	0.61
12:L:74:THR:HG22	12:L:107:PHE:HB2	1.82	0.61
2:2:203:G:O2'	2:2:465:A:N1	2.30	0.61
1:1:381:G:OP1	24:X:18:ARG:NH2	2.33	0.61
1:1:544:C:H2'	1:1:545:U:C4'	2.31	0.61
1:1:1800:C:OP2	4:B:182:ARG:NH1	2.33	0.61
18:R:51:VAL:HG13	18:R:52:PRO:HD3	1.82	0.61
1:1:1852:U:O2	1:1:1890:A:N6	2.34	0.61
2:2:4:U:N3	2:2:6:G:OP1	2.31	0.61
23:W:53:CYS:SG	23:W:57:HIS:ND1	2.69	0.61
1:1:2402:U:O2'	1:1:2403:C:H5''	2.01	0.61
2:2:582:C:OP2	2:2:758:C:N4	2.33	0.61
12:L:58:TYR:O	30:e:13:ARG:NH2	2.34	0.61
34:i:147:GLU:O	34:i:148:LYS:C	2.44	0.61
1:1:571:U:H3'	18:R:80:ARG:HH22	1.66	0.60
1:1:860:U:H2'	1:1:861:A:H8	1.65	0.60
2:2:8:A:N6	34:i:202:GLU:O	2.34	0.60
2:2:1147:C:H1'	39:n:18:ARG:HH21	1.66	0.60
36:k:38:ARG:NH1	36:k:98:GLU:O	2.33	0.60
54:A:110:ASN:HA	54:A:171:VAL:HB	1.83	0.60
1:1:1528:A:N6	1:1:1543:G:O2'	2.33	0.60
6:D:150:THR:HG22	6:D:152:GLU:H	1.67	0.60
39:n:42:GLU:OE2	39:n:45:ARG:NH1	2.34	0.60
1:1:404:A:N6	1:1:421:C:O2'	2.33	0.60
1:1:882:G:H1	1:1:895:U:H4'	1.66	0.60
1:1:1349:C:N4	1:1:1350:C:N4	2.49	0.60
1:1:2831:G:OP2	5:C:59:ARG:NH1	2.34	0.60
2:2:107:G:N7	50:y:10:ARG:NH2	2.48	0.60
2:2:151:A:OP2	2:2:169:C:N4	2.34	0.60
2:2:625:U:H2'	2:2:626:G:H8	1.66	0.60
30:e:32:ILE:O	30:e:32:ILE:HG22	2.00	0.60
39:n:39:PHE:O	39:n:45:ARG:NE	2.35	0.60
2:2:1328:C:O2'	43:r:29:ARG:NH1	2.33	0.60
26:Z:31:ARG:HD2	26:Z:34:HIS:HB2	1.83	0.60
1:1:282:A:N1	1:1:357:C:N4	2.49	0.60
2:2:376:G:H1	2:2:387:U:H3	1.50	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:544:G:OP1	34:i:56:ARG:NH2	2.35	0.60
2:2:811:C:O2'	2:2:901:A:N1	2.34	0.60
2:2:1095:U:OP1	2:2:1108:G:N1	2.33	0.60
32:g:97:LEU:H	32:g:100:MET:HE3	1.66	0.60
33:h:183:ASP:HB3	33:h:202:ILE:HB	1.83	0.60
46:u:45:GLU:HG2	46:u:46:LYS:HG2	1.84	0.60
49:x:33:THR:HG22	49:x:35:SER:H	1.65	0.60
1:1:2684:U:OP2	16:P:51:ARG:NH1	2.34	0.60
2:2:246:A:N1	2:2:278:G:O2'	2.34	0.60
1:1:1175:A:H4'	1:1:1176:U:O5'	2.01	0.60
2:2:403:C:H2'	2:2:404:G:H8	1.66	0.60
40:o:7:ARG:HB2	40:o:101:SER:HB2	1.84	0.60
41:p:28:ASN:O	41:p:57:LYS:NZ	2.34	0.60
1:1:2162:G:OP1	1:1:2171:A:N6	2.35	0.59
1:1:2659:G:N2	1:1:2662:A:OP2	2.34	0.59
24:X:37:ARG:HG2	24:X:48:THR:HG22	1.84	0.59
27:b:54:VAL:HG23	27:b:55:ILE:HG22	1.82	0.59
7:E:35:THR:HB	7:E:155:THR:HB	1.84	0.59
22:V:64:VAL:HG22	22:V:69:GLU:HG3	1.82	0.59
1:1:1528:A:OP2	1:1:1543:G:N2	2.35	0.59
2:2:923:A:N6	2:2:1392:G:O6	2.35	0.59
2:2:993:G:N7	2:2:1213:A:N6	2.51	0.59
29:d:24:THR:HG23	29:d:27:GLY:H	1.68	0.59
54:A:114:GLU:HB2	54:A:204:ALA:HB3	1.83	0.59
1:1:603:A:N7	1:1:654:A:O2'	2.34	0.59
1:1:2333:A:OP1	23:W:77:ARG:NH2	2.36	0.59
5:C:179:ARG:HB3	5:C:188:LEU:HD12	1.84	0.59
2:2:979:C:OP1	2:2:1223:C:N4	2.36	0.59
40:o:8:ILE:HG12	40:o:100:ILE:HG22	1.83	0.59
1:1:1652:A:OP1	14:N:8:ARG:NH2	2.35	0.59
1:1:1744:A:H3'	1:1:1745:A:H8	1.68	0.59
2:2:694:A:N1	2:2:787:A:O2'	2.36	0.59
2:2:1441:A:H2'	2:2:1442:G:C8	2.38	0.59
54:A:124:ALA:HB2	54:A:199:SER:HB3	1.84	0.59
1:1:404:A:N3	1:1:406:G:C6	2.70	0.59
1:1:210:C:OP1	29:d:29:GLN:NE2	2.35	0.59
36:k:10:VAL:HB	36:k:58:HIS:HB3	1.85	0.59
1:1:1077:A:N6	1:1:1088:A:O2'	2.35	0.59
2:2:1103:C:OP1	32:g:95:ARG:NH2	2.34	0.59
2:2:1417:G:O2'	2:2:1483:A:N6	2.35	0.59
13:M:66:ARG:NH1	13:M:104:GLU:OE2	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:404:A:C2	1:1:406:G:C6	2.91	0.58
1:1:2688:G:N1	1:1:2720:U:OP2	2.32	0.58
2:2:1178:G:N2	2:2:1181:G:OP2	2.34	0.58
43:r:25:VAL:HG12	43:r:29:ARG:HD2	1.85	0.58
1:1:102:U:O4	25:Y:2:LYS:N	2.36	0.58
2:2:517:G:N1	2:2:533:A:OP2	2.36	0.58
2:2:1060:U:OP1	44:s:85:ARG:NH1	2.36	0.58
1:1:2058:A:N1	1:1:2505:G:N1	2.51	0.58
5:C:136:ASN:HD21	5:C:139:SER:HB2	1.69	0.58
18:R:38:VAL:O	18:R:54:VAL:CG2	2.40	0.58
33:h:156:ARG:HH11	33:h:193:TYR:HB3	1.68	0.58
1:1:956:G:O6	13:M:14:LYS:NZ	2.36	0.58
1:1:1064:C:O2'	1:1:1075:C:N4	2.36	0.58
2:2:1257:A:H3'	2:2:1258:G:C8	2.39	0.58
1:1:2873:A:O2'	1:1:2874:C:H5'	2.03	0.58
3:3:12:C:O3'	3:3:15:A:N6	2.37	0.58
11:K:105:ARG:NH1	16:P:34:GLU:OE1	2.37	0.58
36:k:10:VAL:HA	36:k:84:VAL:HA	1.86	0.58
1:1:139:U:H5'	1:1:140:C:H5''	1.86	0.58
1:1:785:G:C6	1:1:786:C:C4	2.91	0.58
4:B:66:ASP:OD2	4:B:102:ARG:NH1	2.37	0.58
15:O:43:ASN:HD21	15:O:46:GLU:HG2	1.69	0.58
35:j:97:GLN:HB3	35:j:124:LEU:HB2	1.85	0.58
1:1:1380:G:C6	1:1:1381:G:O6	2.56	0.58
2:2:187:G:O2'	2:2:190:A:N6	2.37	0.58
17:Q:76:TYR:OH	17:Q:92:ARG:NH1	2.37	0.58
1:1:1227:G:OP2	17:Q:16:LYS:NZ	2.37	0.58
2:2:673:A:H4'	36:k:86:ARG:HH12	1.68	0.58
2:2:935:A:H61	37:l:3:ARG:HG3	1.68	0.58
7:E:49:LEU:O	7:E:150:ARG:NH2	2.37	0.58
21:U:86:ARG:NH2	21:U:102:THR:OG1	2.36	0.58
37:l:5:ARG:NH1	37:l:6:VAL:O	2.37	0.58
54:A:225:ASP:HB2	54:A:245:ARG:HB3	1.85	0.58
1:1:396:G:OP2	24:X:10:LYS:NZ	2.36	0.58
2:2:957:U:OP1	49:x:81:ARG:NH1	2.34	0.58
1:1:138:U:H1'	1:1:141:G:H1	1.68	0.58
1:1:957:C:H5'	13:M:75:GLU:HG2	1.86	0.58
1:1:1485:U:H3	1:1:1504:A:H2	1.51	0.58
47:v:17:MET:HG3	47:v:20:SER:HB2	1.86	0.58
1:1:1853:A:N6	1:1:1888:G:O2'	2.36	0.57
1:1:2133:G:N2	1:1:2158:A:OP2	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:29:PRO:HG2	4:B:34:LEU:HD11	1.85	0.57
47:v:61:ILE:HA	47:v:75:LEU:HA	1.86	0.57
1:1:643:A:N1	1:1:2369:A:O2'	2.36	0.57
18:R:43:ASN:HB3	18:R:46:GLU:HB3	1.86	0.57
38:m:5:ASP:OD2	38:m:77:ARG:NH1	2.37	0.57
1:1:2290:G:N2	1:1:2343:U:O2	2.38	0.57
1:1:2314:A:OP1	7:E:88:LYS:NZ	2.37	0.57
39:n:106:ARG:NH1	39:n:107:ASP:O	2.37	0.57
40:o:5:ARG:HB2	40:o:77:VAL:HA	1.86	0.57
1:1:1722:A:N6	1:1:1738:G:O2'	2.37	0.57
6:D:147:LEU:HB2	6:D:183:PHE:HD2	1.69	0.57
2:2:1418:A:N6	2:2:1482:G:O2'	2.38	0.57
6:D:117:ARG:NH1	12:L:1:MET:O	2.37	0.57
16:P:33:VAL:HG22	16:P:38:LYS:HG2	1.87	0.57
33:h:56:VAL:HB	33:h:67:THR:HB	1.87	0.57
1:1:321:U:O3'	6:D:162:ARG:NH1	2.37	0.57
1:1:2127:G:H21	1:1:2173:A:H1'	1.70	0.57
4:B:258:ARG:NH2	4:B:263:THR:OG1	2.37	0.57
14:N:87:PHE:HD1	14:N:90:ARG:HD3	1.70	0.57
18:R:6:GLN:HG2	18:R:11:GLN:HG2	1.87	0.57
21:U:40:ASN:ND2	21:U:64:ALA:O	2.38	0.57
1:1:893:C:H2'	1:1:894:U:C6	2.40	0.57
2:2:409:U:H5'	34:i:25:VAL:CG2	2.35	0.57
52:5:14:G:H22	52:5:48:C:H42	1.53	0.57
1:1:545:U:O2'	1:1:548:G:OP2	2.21	0.57
1:1:1508:A:H4'	1:1:1509:A:H5'	1.87	0.57
8:F:80:THR:HG23	8:F:81:GLU:HG3	1.86	0.57
19:S:25:ARG:NH1	19:S:74:ILE:O	2.36	0.57
1:1:248:G:O5'	1:1:249:C:H5''	2.05	0.56
1:1:1278:C:H2'	1:1:1279:G:H8	1.69	0.56
1:1:1477:A:N6	1:1:1514:G:O2'	2.38	0.56
1:1:1816:C:N4	1:1:2213:U:O4	2.38	0.56
2:2:80:A:OP2	2:2:83:C:N4	2.38	0.56
9:G:129:GLU:HA	9:G:143:ILE:HA	1.86	0.56
12:L:55:MET:O	12:L:60:ARG:NH2	2.38	0.56
1:1:370:G:O2'	1:1:424:G:OP1	2.23	0.56
1:1:408:G:H1	1:1:419:U:H3	1.53	0.56
1:1:1223:G:OP1	18:R:68:ARG:NH2	2.38	0.56
1:1:1392:A:N6	20:T:18:GLU:OE1	2.38	0.56
11:K:108:ARG:NH1	16:P:34:GLU:OE1	2.38	0.56
18:R:74:ILE:HB	18:R:87:GLN:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1380:G:C5	1:1:1381:G:C5	2.93	0.56
1:1:1380:G:O2'	1:1:1381:G:H5'	2.05	0.56
1:1:1935:G:H3'	1:1:1962:5MC:HN41	1.71	0.56
1:1:2830:C:OP2	5:C:59:ARG:NH2	2.38	0.56
2:2:542:G:O3'	34:i:14:ARG:NH2	2.39	0.56
6:D:97:ASN:ND2	6:D:100:MET:SD	2.78	0.56
17:Q:44:GLN:HE21	18:R:77:PHE:HB3	1.70	0.56
39:n:12:ARG:NH1	39:n:107:ASP:OD2	2.37	0.56
54:A:298:LEU:HD23	54:A:344:ILE:HG23	1.86	0.56
2:2:300:A:O2'	2:2:564:C:N3	2.34	0.56
2:2:727:G:N2	2:2:730:G:OP2	2.36	0.56
34:i:57:GLU:HG2	34:i:199:LEU:HB2	1.86	0.56
39:n:24:GLY:H	39:n:61:LEU:HA	1.70	0.56
2:2:615:G:C6	2:2:626:G:C6	2.94	0.56
7:E:119:ALA:O	7:E:167:ARG:NH2	2.35	0.56
36:k:63:ASN:ND2	36:k:96:VAL:O	2.38	0.56
54:A:138:TYR:OH	54:A:142:ARG:NH2	2.38	0.56
1:1:335:C:O2	21:U:68:SER:OG	2.23	0.56
3:3:86:G:H1	3:3:90:C:H42	1.53	0.56
1:1:796:C:OP1	6:D:57:LYS:NZ	2.39	0.56
1:1:1666:G:N3	11:K:3:GLN:NE2	2.54	0.56
4:B:161:TYR:HB3	4:B:194:GLU:HG2	1.88	0.56
1:1:1450:G:H1	1:1:1461:C:H42	1.51	0.56
1:1:1722:A:OP2	1:1:1738:G:N2	2.39	0.56
1:1:2305:U:H5''	7:E:131:GLY:HA3	1.86	0.56
1:1:2818:U:OP2	14:N:42:LYS:NZ	2.36	0.56
26:Z:6:LYS:HB2	26:Z:58:GLU:HB3	1.87	0.56
1:1:1009:A:OP1	10:J:39:LYS:NZ	2.39	0.56
1:1:1188:U:H2'	1:1:1189:A:H8	1.71	0.56
1:1:2087:G:H2'	1:1:2088:A:H8	1.71	0.56
1:1:2141:G:H1	1:1:2150:C:H42	1.54	0.56
2:2:1275:A:H3'	2:2:1276:G:H8	1.71	0.56
54:A:244:ILE:HG21	54:A:267:LYS:HA	1.87	0.56
1:1:1869:G:O2'	1:1:1872:A:N6	2.39	0.55
2:2:944:G:N1	2:2:1338:G:OP2	2.39	0.55
2:2:1414:U:H3	2:2:1486:G:H1	1.52	0.55
34:i:121:LYS:C	34:i:146:ARG:HD3	2.29	0.55
42:q:27:CYS:HB2	42:q:30:LYS:HD3	1.88	0.55
1:1:1313:U:O3'	1:1:1332:G:H5''	2.07	0.55
1:1:2295:C:OP1	15:O:10:ARG:NH1	2.37	0.55
2:2:297:G:N2	2:2:300:A:OP2	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:1314:C:N4	49:x:2:PRO:O	2.39	0.55
22:V:72:VAL:HG12	22:V:93:ARG:HA	1.87	0.55
1:1:1067:A:N3	1:1:1068:G:N1	2.52	0.55
33:h:46:GLU:HG2	33:h:87:LEU:HD21	1.88	0.55
33:h:140:ASN:OD1	33:h:143:ARG:NH2	2.38	0.55
15:O:33:ARG:O	15:O:65:THR:OG1	2.25	0.55
48:w:48:ARG:HH22	48:w:50:LYS:HD3	1.71	0.55
1:1:914:G:H5'	1:1:915:C:OP2	2.07	0.55
1:1:1026:G:H2'	1:1:1027:A:H8	1.71	0.55
1:1:2680:U:H5'	5:C:194:PRO:HA	1.87	0.55
2:2:1432:G:O2'	2:2:1467:C:N4	2.39	0.55
2:2:1450:U:O2'	2:2:1453:G:O6	2.25	0.55
7:E:4:LEU:HA	7:E:7:TYR:HB3	1.89	0.55
19:S:73:LYS:HB2	19:S:106:VAL:HB	1.88	0.55
24:X:33:LEU:HD23	24:X:50:ARG:HG2	1.87	0.55
1:1:404:A:HO2'	1:1:405:U:P	2.18	0.55
1:1:714:U:OP2	45:t:88:ARG:NH1	2.33	0.55
1:1:807:U:H2'	1:1:808:G:H8	1.72	0.55
2:2:62:U:H3	2:2:105:G:H1	1.54	0.55
2:2:1156:G:O2'	2:2:1179:A:N6	2.39	0.55
2:2:1176:A:H2'	2:2:1177:G:C8	2.42	0.55
2:2:1268:G:N3	2:2:1326:U:O2'	2.40	0.55
10:J:72:LYS:HE3	10:J:74:TYR:HE1	1.72	0.55
38:m:64:LYS:HG2	38:m:71:VAL:HG21	1.88	0.55
1:1:688:U:H2'	1:1:689:A:H8	1.72	0.55
1:1:1124:G:N3	31:f:37:GLN:NE2	2.55	0.55
1:1:2419:U:OP1	30:e:41:LYS:NZ	2.37	0.55
34:i:60:LYS:NZ	34:i:194:ASP:O	2.40	0.55
1:1:83:A:N6	1:1:101:A:O2'	2.40	0.55
1:1:1412:U:H2'	1:1:1413:A:H8	1.71	0.55
1:1:2391:G:O2'	1:1:2424:C:N4	2.40	0.55
2:2:160:A:H1'	2:2:344:A:C5	2.42	0.55
2:2:611:C:H42	2:2:629:A:H61	1.52	0.55
2:2:1516:2MG:N2	2:2:1519:MA6:OP2	2.38	0.55
8:F:98:VAL:HG22	8:F:103:ILE:HG12	1.88	0.55
37:l:15:ASP:OD1	37:l:20:SER:N	2.40	0.55
1:1:271:G:OP1	1:1:367:G:N2	2.39	0.55
1:1:1996:C:OP2	11:K:31:ARG:NH2	2.40	0.55
2:2:559:A:H4'	2:2:560:A:H3'	1.89	0.55
2:2:1380:U:H4'	2:2:1381:U:H5'	1.88	0.55
3:3:7:G:O2'	15:O:38:GLN:NE2	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:F:17:VAL:HG22	8:F:26:ILE:HG12	1.88	0.55
1:1:1349:C:H42	1:1:1382:G:H1	1.54	0.54
1:1:1482:G:O2'	1:1:1509:A:N6	2.38	0.54
1:1:1509:A:H1'	1:1:1510:G:C8	2.42	0.54
1:1:1736:U:H2'	1:1:1737:G:O4'	2.06	0.54
1:1:1844:C:O3'	4:B:256:LYS:NZ	2.40	0.54
2:2:131:A:O2'	2:2:262:A:N3	2.39	0.54
43:r:24:GLY:HA3	43:r:65:VAL:HG12	1.88	0.54
1:1:1040:A:O2'	1:1:1116:G:N2	2.38	0.54
1:1:1467:U:H5	1:1:1546:G:H2'	1.72	0.54
31:f:2:LYS:NZ	31:f:32:LYS:O	2.39	0.54
37:l:40:GLU:OE1	39:n:41:ARG:NH1	2.39	0.54
1:1:706:A:OP1	4:B:7:LYS:NZ	2.34	0.54
1:1:881:G:N2	1:1:882:G:O6	2.40	0.54
1:1:1693:U:O2	4:B:14:ARG:NH1	2.40	0.54
1:1:1716:U:H2'	1:1:1717:A:H8	1.73	0.54
1:1:1901:A:OP2	4:B:253:LYS:NZ	2.37	0.54
2:2:595:A:O2'	2:2:640:A:N6	2.39	0.54
5:C:181:ASP:OD2	5:C:184:ARG:NH2	2.39	0.54
39:n:6:TYR:HB2	39:n:21:ILE:HB	1.89	0.54
39:n:36:GLU:OE1	39:n:45:ARG:NH1	2.40	0.54
1:1:2261:C:OP1	23:W:19:LYS:NZ	2.33	0.54
2:2:181:A:O2'	2:2:194:C:N4	2.40	0.54
10:J:98:GLU:OE2	10:J:126:ALA:N	2.37	0.54
15:O:40:ILE:HG12	15:O:47:VAL:HG12	1.89	0.54
1:1:249:C:O2	30:e:12:LYS:NZ	2.40	0.54
1:1:1857:G:N2	1:1:1884:G:O2'	2.38	0.54
2:2:409:U:H5'	34:i:25:VAL:HG21	1.89	0.54
21:U:46:GLN:NE2	21:U:54:GLN:OE1	2.40	0.54
1:1:532:A:OP1	1:1:561:G:N2	2.39	0.54
1:1:768:G:N2	1:1:1379:U:H1'	2.23	0.54
2:2:197:A:N1	2:2:220:G:O2'	2.37	0.54
14:N:98:LEU:O	14:N:112:TYR:N	2.40	0.54
1:1:981:A:OP2	1:1:982:C:N4	2.37	0.54
1:1:2137:U:N3	1:1:2155:U:O2	2.41	0.54
2:2:1158:C:O2'	2:2:1160:G:OP1	2.26	0.54
3:3:31:C:H2'	3:3:32:U:H6	1.72	0.54
34:i:75:TYR:OH	34:i:97:ARG:NH1	2.41	0.54
24:X:17:ASN:OD1	24:X:27:ARG:NH1	2.41	0.54
44:s:41:ARG:NH2	49:x:6:LYS:O	2.41	0.54
52:5:8:4SU:HN3	52:5:23:C:H41	1.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1069:A:O2'	1:1:1074:G:N2	2.41	0.54
2:2:409:U:H2'	2:2:410:G:O4'	2.08	0.54
2:2:615:G:C2	2:2:626:G:C5	2.96	0.54
2:2:993:G:O2'	2:2:994:A:N7	2.40	0.54
8:F:10:VAL:HA	8:F:49:THR:HA	1.90	0.54
54:A:102:PRO:O	54:A:352:GLN:NE2	2.41	0.54
1:1:954:G:H1	1:1:963:U:H3	1.56	0.54
1:1:1047:G:HO2'	1:1:1110:G:H1	1.54	0.54
2:2:1316:G:N1	2:2:1319:A:OP2	2.40	0.54
33:h:58:GLU:HG2	33:h:60:PRO:HD3	1.90	0.54
1:1:1651:G:N7	14:N:11:ASN:ND2	2.54	0.53
1:1:2175:C:H2'	1:1:2176:A:O4'	2.09	0.53
2:2:54:C:H2'	2:2:352:C:H41	1.72	0.53
2:2:210:C:H4'	2:2:211:G:H5''	1.90	0.53
2:2:1166:G:N2	2:2:1169:A:OP1	2.40	0.53
8:F:2:SER:OG	8:F:6:LYS:NZ	2.42	0.53
54:A:242:SER:O	54:A:266:ASN:ND2	2.40	0.53
1:1:238:C:H42	1:1:259:G:H1	1.56	0.53
2:2:1382:C:O2'	37:l:79:ARG:NH1	2.42	0.53
3:3:36:C:N4	3:3:49:C:O2	2.39	0.53
40:o:14:ASP:HB3	40:o:17:LEU:HB3	1.90	0.53
1:1:411:G:OP2	1:1:2406:A:O2'	2.26	0.53
1:1:568:U:H1'	1:1:2030:6MZ:H9C1	1.91	0.53
2:2:83:C:HO2'	2:2:86:G:H1	1.51	0.53
2:2:159:G:O2'	2:2:162:A:N6	2.42	0.53
7:E:99:PHE:HA	7:E:102:ARG:HE	1.74	0.53
25:Y:49:ASP:OD1	25:Y:52:ARG:NH2	2.42	0.53
35:j:105:ILE:O	35:j:112:ARG:NH1	2.41	0.53
46:u:12:LYS:HG2	46:u:13:LYS:HG2	1.90	0.53
1:1:833:A:N6	1:1:944:C:O2	2.40	0.53
1:1:1869:G:N2	1:1:1871:A:O2'	2.41	0.53
1:1:2299:U:OP1	7:E:72:LYS:NZ	2.37	0.53
1:1:2749:A:OP1	8:F:2:SER:N	2.41	0.53
2:2:579:A:H4'	45:t:54:ARG:HH22	1.73	0.53
1:1:1415:U:O4'	1:1:1588:G:N2	2.41	0.53
1:1:1862:G:H1	1:1:1880:U:H3	1.56	0.53
1:1:1964:G:O2'	1:1:1967:C:OP2	2.27	0.53
2:2:266:G:H3'	47:v:69:LYS:HB2	1.89	0.53
2:2:322:C:O2	2:2:332:G:N2	2.41	0.53
2:2:1304:G:H21	2:2:1333:A:H62	1.56	0.53
1:1:61:C:H4'	25:Y:40:SER:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:447:A:OP1	17:Q:5:LYS:NZ	2.40	0.53
1:1:567:U:OP2	12:L:29:LYS:NZ	2.41	0.53
1:1:1534:U:O2'	1:1:1537:G:N1	2.34	0.53
1:1:2266:A:N6	1:1:2273:A:OP2	2.41	0.53
1:1:2626:C:H2'	1:1:2627:G:C8	2.43	0.53
2:2:832:G:H1	2:2:854:U:H3	1.54	0.53
2:2:1311:A:H2	2:2:1326:U:H3	1.57	0.53
4:B:93:LEU:HD11	4:B:101:ARG:HB3	1.89	0.53
1:1:500:G:N1	1:1:503:A:OP2	2.41	0.53
1:1:1580:A:H3'	1:1:1581:G:H8	1.73	0.53
1:1:2403:C:H42	1:1:2414:G:H1	1.57	0.53
4:B:2:ALA:N	4:B:20:VAL:O	2.42	0.53
13:M:14:LYS:O	13:M:71:LYS:NZ	2.32	0.53
1:1:720:U:H2'	1:1:721:A:C8	2.44	0.53
1:1:1446:C:HO2'	1:1:1545:A:HO2'	1.56	0.53
1:1:1939:5MU:OP1	1:1:2604:U:O2'	2.27	0.53
2:2:129:A:H2	2:2:232:G:H22	1.57	0.53
4:B:258:ARG:NH1	4:B:264:ASP:OD1	2.35	0.53
16:P:88:ARG:NH2	16:P:110:ILE:O	2.38	0.53
21:U:72:ILE:HD12	21:U:83:VAL:HG21	1.91	0.53
1:1:142:A:N3	20:T:1:MET:N	2.56	0.53
1:1:673:C:OP1	6:D:49:ARG:NH2	2.38	0.53
1:1:1476:U:OP2	1:1:1514:G:N1	2.38	0.53
2:2:974:A:OP1	44:s:69:ARG:NH1	2.41	0.53
4:B:50:THR:OG1	4:B:51:THR:N	2.42	0.53
1:1:51:G:O2'	1:1:119:A:N1	2.42	0.53
1:1:463:G:N2	1:1:466:A:OP2	2.41	0.53
1:1:857:G:N2	1:1:921:C:O2	2.42	0.53
1:1:2639:A:O3'	10:J:96:ARG:NH2	2.42	0.53
2:2:1027:C:H42	2:2:1033:G:H1	1.57	0.53
2:2:1152:A:OP1	40:o:70:HIS:ND1	2.42	0.53
2:2:1213:A:O2'	2:2:1215:G:N7	2.36	0.53
3:3:63:C:H2'	3:3:64:G:H8	1.74	0.53
5:C:152:PRO:O	5:C:154:LYS:N	2.42	0.53
33:h:136:ARG:O	33:h:140:ASN:ND2	2.43	0.53
1:1:1251:C:OP2	17:Q:6:ARG:NH2	2.35	0.52
1:1:2473:U:OP1	1:1:2529:G:N2	2.42	0.52
39:n:119:ARG:HG3	39:n:125:PRO:HG3	1.90	0.52
54:A:223:ARG:HB3	54:A:247:THR:HB	1.91	0.52
1:1:277:G:N2	1:1:359:G:O5'	2.41	0.52
1:1:601:C:O2'	6:D:99:LYS:NZ	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2638:G:O2'	1:1:2775:G:N2	2.41	0.52
2:2:720:C:O3'	48:w:52:GLN:NE2	2.42	0.52
4:B:69:ARG:NH2	4:B:127:GLY:O	2.39	0.52
52:5:7:U:H2'	52:5:12:C:H41	1.74	0.52
2:2:137:U:H2'	2:2:138:G:H8	1.74	0.52
2:2:346:G:OP1	16:P:39:ARG:NH2	2.43	0.52
8:F:99:LYS:HG2	8:F:104:ASN:HD21	1.74	0.52
1:1:994:C:O2'	1:1:996:A:OP1	2.24	0.52
1:1:1350:C:C4	1:1:1351:C:C4	2.97	0.52
2:2:55:A:N6	2:2:356:A:N1	2.58	0.52
9:G:116:ARG:NE	9:G:133:GLN:OE1	2.42	0.52
20:T:65:GLY:N	20:T:79:ASP:OD1	2.42	0.52
21:U:74:ASN:ND2	21:U:81:ASP:OD2	2.40	0.52
1:1:1426:G:O2'	1:1:1572:A:N6	2.43	0.52
1:1:1863:G:H1	1:1:1879:C:H42	1.57	0.52
1:1:2208:C:H2'	1:1:2209:G:C8	2.45	0.52
2:2:1086:U:H2'	2:2:1087:G:H8	1.73	0.52
1:1:879:G:N2	1:1:899:A:O2'	2.42	0.52
1:1:1857:G:O2'	1:1:1884:G:N2	2.41	0.52
1:1:1994:C:H2'	1:1:1995:U:H6	1.75	0.52
1:1:2252:G:O6	13:M:81:ARG:NH2	2.42	0.52
2:2:1526:G:N7	51:z:40:LYS:NZ	2.49	0.52
7:E:68:THR:N	7:E:86:GLY:O	2.43	0.52
2:2:832:G:OP2	32:g:21:ARG:NH2	2.43	0.52
2:2:1004:A:N6	2:2:1024:G:O2'	2.43	0.52
2:2:1181:G:H4'	2:2:1182:G:H5'	1.92	0.52
2:2:1350:A:OP2	39:n:120:LYS:NZ	2.38	0.52
34:i:121:LYS:O	34:i:146:ARG:CD	2.47	0.52
52:5:18:G:N2	52:5:58:A:OP2	2.42	0.52
1:1:320:A:O2'	1:1:322:A:OP2	2.28	0.52
1:1:1064:C:N4	1:1:1070:A:OP2	2.34	0.52
2:2:309:A:O2'	2:2:607:A:N1	2.42	0.52
2:2:842:U:O2'	2:2:845:A:OP2	2.24	0.52
2:2:977:A:O3'	2:2:980:C:N4	2.43	0.52
15:O:71:ALA:HB1	15:O:106:LEU:HB2	1.91	0.52
34:i:145:ILE:C	34:i:147:GLU:H	2.17	0.52
39:n:19:VAL:HA	39:n:65:ILE:HG12	1.90	0.52
39:n:116:VAL:HG21	40:o:62:ARG:HB2	1.91	0.52
1:1:747:5MU:O2'	19:S:92:ARG:NH1	2.43	0.52
1:1:1089:A:H2	1:1:1090:A:H62	1.58	0.52
2:2:662:U:O2'	2:2:836:G:OP1	2.28	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:1129:C:O4'	2:2:1146:A:N6	2.41	0.52
2:2:1258:G:H2'	2:2:1259:C:C6	2.45	0.52
7:E:47:LYS:HG2	7:E:51:ASP:HB2	1.91	0.52
44:s:24:ARG:HG2	44:s:51:LEU:HD13	1.92	0.52
1:1:72:U:OP2	25:Y:54:LYS:NZ	2.43	0.52
1:1:2646:C:OP2	1:1:2732:G:O2'	2.25	0.52
1:1:2743:U:OP2	1:1:2755:C:N4	2.43	0.52
2:2:561:U:O2	42:q:15:LYS:NZ	2.41	0.52
2:2:911:U:OP2	42:q:94:ARG:NH2	2.42	0.52
2:2:1376:U:H2'	2:2:1377:A:H8	1.75	0.52
1:1:1306:C:H41	1:1:1606:C:H2'	1.74	0.51
1:1:2008:C:H2'	1:1:2009:A:H8	1.74	0.51
2:2:212:G:N2	2:2:213:G:O6	2.44	0.51
6:D:168:ASP:OD2	6:D:170:ARG:NE	2.43	0.51
1:1:275:C:O2'	1:1:362:A:N6	2.43	0.51
1:1:353:C:H3'	1:1:354:A:H8	1.75	0.51
1:1:376:G:H2'	1:1:377:G:H8	1.75	0.51
1:1:2233:U:H2'	1:1:2234:G:H8	1.75	0.51
1:1:2357:G:N2	1:1:2360:G:OP2	2.35	0.51
2:2:1234:C:H1'	2:2:1364:U:H6	1.76	0.51
5:C:25:THR:HG21	5:C:193:VAL:HG22	1.90	0.51
7:E:134:GLU:HB3	7:E:137:ILE:HG23	1.91	0.51
18:R:4:VAL:O	18:R:39:LEU:N	2.38	0.51
26:Z:15:GLY:O	26:Z:16:ARG:NH1	2.42	0.51
32:g:130:THR:HG22	32:g:132:LYS:HG2	1.92	0.51
41:p:22:HIS:HB2	41:p:33:THR:HB	1.92	0.51
1:1:1865:U:O2'	1:1:1875:G:N2	2.37	0.51
1:1:1997:C:H2'	1:1:1998:A:H8	1.75	0.51
1:1:2527:C:H5'	31:f:34:LYS:HD3	1.91	0.51
1:1:2595:G:N2	1:1:2598:A:OP2	2.35	0.51
32:g:66:LYS:O	32:g:159:ASP:N	2.43	0.51
35:j:38:VAL:HA	35:j:48:PHE:HA	1.93	0.51
49:x:11:ILE:HG13	49:x:38:SER:HB3	1.91	0.51
1:1:318:C:H2'	1:1:319:G:H8	1.74	0.51
1:1:2339:C:H6	1:1:2339:C:H5''	1.76	0.51
1:1:2484:G:OP1	13:M:44:ARG:NH1	2.44	0.51
1:1:2756:U:H5''	31:f:19:ARG:HG2	1.92	0.51
2:2:70:U:O2	2:2:71:A:N6	2.39	0.51
2:2:255:G:H4'	47:v:19:LYS:HB2	1.93	0.51
4:B:180:GLU:OE2	4:B:270:ARG:NH1	2.43	0.51
6:D:58:LYS:NZ	6:D:70:SER:O	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:c:25:LYS:NZ	28:c:32:GLU:O	2.43	0.51
1:1:542:C:H2'	1:1:543:G:C8	2.46	0.51
1:1:2133:G:O3'	1:1:2158:A:N6	2.44	0.51
2:2:1302:C:H5'	43:r:17:ILE:HG23	1.91	0.51
2:2:1449:C:H2'	2:2:1450:U:O4'	2.11	0.51
22:V:58:SER:O	22:V:73:LYS:NZ	2.42	0.51
31:f:2:LYS:HB2	31:f:35:GLN:HG2	1.92	0.51
37:l:24:ALA:O	37:l:28:ASN:ND2	2.43	0.51
37:l:140:ASP:OD1	37:l:143:ARG:NH1	2.43	0.51
2:2:757:U:HO2'	2:2:879:C:HO2'	1.49	0.51
2:2:1006:G:O6	2:2:1024:G:N2	2.43	0.51
2:2:1279:G:H4'	2:2:1281:C:H5	1.76	0.51
21:U:4:LYS:O	21:U:94:ARG:NH2	2.40	0.51
24:X:2:SER:OG	24:X:3:ARG:N	2.35	0.51
34:i:188:ARG:NH2	34:i:195:ILE:O	2.42	0.51
1:1:880:G:H22	1:1:898:C:H1'	1.74	0.51
1:1:1045:C:H4'	1:1:1046:A:H5''	1.92	0.51
1:1:1649:G:O2'	14:N:106:ASP:OD2	2.27	0.51
2:2:1458:G:OP1	50:y:30:THR:OG1	2.29	0.51
1:1:783:A:N3	1:1:783:A:C2'	2.73	0.51
1:1:2271:G:H5'	23:W:20:ARG:HG2	1.93	0.51
2:2:49:U:H3	2:2:362:G:H1'	1.76	0.51
2:2:127:G:O2'	47:v:6:ARG:NH1	2.44	0.51
11:K:21:CYS:HA	11:K:41:ILE:HG22	1.92	0.51
13:M:50:ARG:HD3	13:M:65:ILE:HD11	1.93	0.51
32:g:24:ASN:ND2	32:g:192:ASP:OD1	2.43	0.51
1:1:210:C:OP1	29:d:25:LYS:NZ	2.35	0.51
2:2:564:C:OP1	42:q:12:ARG:NH1	2.44	0.51
2:2:1111:A:N6	33:h:176:HIS:O	2.43	0.51
2:2:1233:G:O5'	39:n:119:ARG:NH1	2.44	0.51
7:E:58:ALA:HB2	7:E:65:PRO:HD3	1.92	0.51
22:V:86:LEU:HD13	22:V:89:ILE:HD11	1.93	0.51
33:h:47:LEU:HB3	33:h:50:ALA:HB3	1.93	0.51
37:l:118:LEU:O	37:l:122:ASN:ND2	2.44	0.51
1:1:518:G:O5'	19:S:18:ARG:NH1	2.44	0.51
1:1:2298:A:H61	1:1:2318:G:H1'	1.76	0.51
2:2:126:G:OP1	2:2:605:U:O2'	2.24	0.51
2:2:1114:C:H42	2:2:1186:G:H1	1.57	0.51
2:2:1377:A:OP1	37:l:92:ARG:NH1	2.38	0.51
4:B:168:ASP:N	4:B:168:ASP:OD1	2.43	0.51
1:1:192:C:O2'	1:1:802:A:N3	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1263:U:OP1	27:b:13:ARG:NH1	2.44	0.50
1:1:1274:A:OP1	1:1:1646:C:N4	2.42	0.50
2:2:35:G:O2'	42:q:115:SER:O	2.29	0.50
54:A:110:ASN:O	54:A:208:GLU:N	2.44	0.50
1:1:220:G:O2'	1:1:233:A:N3	2.35	0.50
1:1:1095:A:O2'	1:1:1096:A:O4'	2.30	0.50
2:2:373:A:H61	2:2:391:G:H1'	1.76	0.50
2:2:418:C:H42	2:2:425:G:H1	1.59	0.50
32:g:54:LEU:HD23	32:g:57:LEU:HD12	1.92	0.50
51:z:42:THR:OG1	51:z:43:THR:N	2.45	0.50
1:1:250:G:O2'	12:L:59:ARG:NH2	2.44	0.50
1:1:1477:A:OP2	1:1:1514:G:N2	2.44	0.50
2:2:958:A:N3	2:2:985:C:O2'	2.40	0.50
7:E:47:LYS:O	7:E:51:ASP:N	2.42	0.50
9:G:132:PHE:HB2	9:G:140:ALA:HB3	1.92	0.50
43:r:98:ARG:HB2	43:r:100:GLN:HE22	1.77	0.50
54:A:152:SER:O	54:A:160:LYS:N	2.44	0.50
2:2:324:G:N1	2:2:327:A:OP2	2.44	0.50
2:2:977:A:N6	2:2:1224:U:O4'	2.43	0.50
2:2:1414:U:H2'	2:2:1415:G:C8	2.46	0.50
4:B:142:HIS:ND1	4:B:193:GLY:O	2.36	0.50
7:E:30:ARG:H	7:E:159:THR:HG1	1.59	0.50
32:g:19:GLN:HB2	32:g:22:TYR:HD2	1.77	0.50
1:1:2451:A:N1	53:H:78:PHE:HE1	2.10	0.50
2:2:1009:U:H2'	2:2:1020:G:H22	1.77	0.50
13:M:47:GLU:OE2	13:M:50:ARG:NH1	2.44	0.50
35:j:15:LEU:HA	35:j:37:THR:HG22	1.94	0.50
37:l:79:ARG:H	37:l:85:TYR:HB2	1.76	0.50
1:1:611:C:H2'	1:1:612:G:C8	2.47	0.50
1:1:901:C:H2'	1:1:902:C:O4'	2.11	0.50
1:1:1469:A:OP2	1:1:1522:A:N6	2.45	0.50
2:2:413:G:N2	2:2:428:G:O3'	2.44	0.50
2:2:802:A:H3'	2:2:803:G:H8	1.77	0.50
6:D:148:ILE:HB	6:D:169:VAL:HG22	1.93	0.50
34:i:150:LYS:NZ	34:i:176:GLY:O	2.39	0.50
38:m:115:ALA:O	38:m:119:ALA:N	2.44	0.50
39:n:129:LYS:NZ	52:5:33:U:OP1	2.41	0.50
54:A:217:ILE:HD11	54:A:271:LEU:HD11	1.94	0.50
1:1:355:U:H2'	1:1:356:G:C8	2.46	0.50
1:1:893:C:O2'	1:1:894:U:C6	2.65	0.50
1:1:1379:U:H2'	1:1:1379:U:O2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1657:U:OP1	5:C:141:ARG:N	2.44	0.50
1:1:2258:C:O2'	1:1:2427:C:OP2	2.28	0.50
1:1:2585:U:C2	53:H:79:PHE:HE1	2.30	0.50
2:2:1274:A:O2'	2:2:1275:A:O4'	2.28	0.50
2:2:1346:A:OP1	39:n:122:ARG:NH1	2.35	0.50
20:T:25:GLU:HG3	20:T:26:LYS:HG3	1.93	0.50
49:x:25:SER:OG	49:x:27:ASP:OD1	2.29	0.50
1:1:445:C:OP1	17:Q:2:ALA:N	2.45	0.50
1:1:2559:C:H2'	1:1:2560:A:H8	1.77	0.50
1:1:2850:A:OP2	1:1:2866:U:N3	2.43	0.50
2:2:890:G:O2'	2:2:906:A:N6	2.45	0.50
12:L:133:ALA:O	12:L:137:ALA:N	2.41	0.50
15:O:94:ARG:HD2	15:O:97:PHE:H	1.76	0.50
43:r:34:LEU:HD22	43:r:39:ILE:HB	1.92	0.50
1:1:1350:C:N3	1:1:1351:C:C4	2.80	0.50
2:2:496:A:H2'	2:2:496:A:N3	2.27	0.50
2:2:1123:U:H4'	40:o:39:PRO:HD2	1.94	0.50
2:2:1211:U:H4'	2:2:1212:U:H5''	1.93	0.50
6:D:106:LYS:HG3	6:D:200:LEU:HD23	1.93	0.50
6:D:149:ILE:HD11	6:D:188:MET:HG2	1.93	0.50
13:M:17:ASN:O	13:M:38:ARG:NH1	2.40	0.50
19:S:10:ALA:HB3	19:S:101:SER:HB2	1.94	0.50
47:v:12:VAL:HB	47:v:57:ASP:H	1.76	0.50
1:1:1350:C:N4	1:1:1351:C:N4	2.60	0.49
1:1:2691:C:HO2'	1:1:2871:U:HO2'	1.55	0.49
6:D:125:SER:O	6:D:137:LYS:NZ	2.45	0.49
9:G:47:PHE:HA	9:G:51:ARG:HB2	1.94	0.49
41:p:67:ALA:HB2	41:p:96:THR:HG23	1.94	0.49
1:1:1807:G:N2	1:1:1810:A:OP2	2.42	0.49
1:1:1947:C:HO2'	2:2:1483:A:HO2'	1.61	0.49
2:2:195:A:H2'	2:2:196:A:C4	2.47	0.49
2:2:369:G:OP2	2:2:388:G:N1	2.44	0.49
2:2:1229:A:OP2	43:r:113:ARG:NH1	2.45	0.49
2:2:1310:G:OP2	43:r:87:ARG:NH1	2.38	0.49
4:B:155:ALA:HB2	4:B:162:VAL:HG23	1.94	0.49
6:D:168:ASP:OD2	6:D:170:ARG:CG	2.60	0.49
6:D:176:ASP:OD1	6:D:176:ASP:N	2.45	0.49
13:M:47:GLU:OE2	13:M:51:ARG:NE	2.45	0.49
35:j:16:ILE:HD12	35:j:36:LEU:HG	1.95	0.49
42:q:33:VAL:N	42:q:56:ARG:O	2.40	0.49
44:s:2:ALA:N	44:s:67:THR:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:236:C:H2'	1:1:237:C:C6	2.47	0.49
1:1:793:A:OP2	1:1:2071:A:O2'	2.27	0.49
1:1:2718:G:O2'	1:1:2847:U:OP1	2.27	0.49
2:2:1328:C:H5''	43:r:28:THR:HG21	1.94	0.49
15:O:111:ARG:NH2	15:O:117:PHE:OXT	2.45	0.49
36:k:100:SER:HB2	36:k:103:VAL:HG23	1.95	0.49
54:A:185:GLN:HG2	54:A:198:THR:HG23	1.94	0.49
1:1:277:G:H1'	1:1:361:G:C6	2.47	0.49
1:1:768:G:H22	1:1:1379:U:H1'	1.77	0.49
1:1:1266:G:O2'	1:1:2012:G:O6	2.28	0.49
1:1:2788:C:O2'	1:1:2809:A:N3	2.43	0.49
2:2:451:A:H61	2:2:481:G:H5'	1.78	0.49
2:2:769:G:N2	2:2:810:C:O2	2.37	0.49
2:2:981:U:O4	2:2:1223:C:N4	2.42	0.49
15:O:68:LYS:HA	15:O:102:ARG:HG2	1.94	0.49
16:P:106:LYS:HD3	16:P:109:ARG:HH22	1.77	0.49
54:A:255:VAL:HG11	54:A:273:VAL:HB	1.94	0.49
1:1:404:A:C4	1:1:406:G:C2	3.01	0.49
1:1:1071:G:H1'	1:1:1089:A:H2'	1.94	0.49
1:1:1754:A:N1	1:1:2716:C:O2'	2.43	0.49
1:1:2333:A:H5'	1:1:2335:A:H1'	1.93	0.49
2:2:80:A:N6	2:2:87:C:N3	2.61	0.49
2:2:157:U:H2'	2:2:158:G:H8	1.77	0.49
19:S:12:SER:OG	19:S:13:SER:N	2.46	0.49
33:h:20:SER:HG	33:h:40:ARG:HH21	1.60	0.49
1:1:307:G:N1	1:1:310:A:OP2	2.44	0.49
1:1:482:A:N6	1:1:506:G:O2'	2.46	0.49
1:1:1788:C:OP1	4:B:221:ARG:NH2	2.44	0.49
1:1:2376:A:N3	15:O:111:ARG:NH1	2.60	0.49
2:2:412:A:H2'	2:2:414:A:H5''	1.95	0.49
2:2:563:A:O2'	2:2:566:G:O3'	2.31	0.49
22:V:51:GLN:HG2	22:V:86:LEU:HD11	1.94	0.49
32:g:94:HIS:ND1	32:g:146:ASN:OD1	2.45	0.49
32:g:119:THR:O	32:g:124:GLY:N	2.42	0.49
39:n:24:GLY:H	39:n:62:ASP:H	1.61	0.49
1:1:1278:C:H2'	1:1:1279:G:C8	2.48	0.49
1:1:2303:G:HO2'	7:E:121:SER:HG	1.60	0.49
2:2:521:G:H4'	42:q:70:GLU:HG2	1.94	0.49
2:2:581:G:N2	2:2:760:G:N7	2.60	0.49
2:2:1066:C:N4	2:2:1191:A:N7	2.61	0.49
5:C:33:ARG:NH1	5:C:53:GLY:O	2.38	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:T:28:ASN:ND2	20:T:88:LYS:O	2.45	0.49
42:q:54:ARG:HD2	42:q:62:GLU:HG3	1.94	0.49
50:y:35:VAL:HG11	50:y:79:LEU:HD13	1.93	0.49
52:5:6:G:O2'	52:5:49:G:O4'	2.30	0.49
52:5:76:A:O2'	53:H:79:PHE:C	2.54	0.49
1:1:1636:U:H2'	1:1:1637:A:H8	1.78	0.49
2:2:1218:C:H2'	2:2:1219:A:C8	2.47	0.49
9:G:47:PHE:O	9:G:52:ALA:N	2.40	0.49
34:i:62:ARG:NH1	34:i:69:GLU:OE1	2.46	0.49
1:1:319:G:H1	1:1:323:C:H42	1.60	0.49
1:1:1994:C:H2'	1:1:1995:U:C6	2.48	0.49
1:1:2543:G:H2'	1:1:2544:G:C8	2.47	0.49
2:2:776:G:N1	2:2:802:A:OP1	2.45	0.49
35:j:88:VAL:HG12	35:j:93:ARG:HA	1.95	0.49
1:1:531:C:H4'	1:1:532:A:H5''	1.95	0.49
1:1:1365:A:OP2	24:X:2:SER:OG	2.30	0.49
1:1:2171:A:O2'	1:1:2174:C:OP2	2.30	0.49
2:2:83:C:O2'	2:2:86:G:N1	2.44	0.49
2:2:104:G:OP2	50:y:13:GLN:NE2	2.43	0.49
2:2:203:G:N2	2:2:204:G:O6	2.35	0.49
8:F:123:ALA:HA	8:F:133:LEU:HA	1.94	0.49
17:Q:24:TYR:HB2	17:Q:29:SER:HB3	1.95	0.49
40:o:40:ILE:O	40:o:73:LEU:N	2.44	0.49
51:z:31:GLU:OE2	51:z:34:ARG:NH2	2.41	0.49
1:1:279:A:OP2	1:1:361:G:N2	2.43	0.48
1:1:404:A:C2	1:1:406:G:N1	2.81	0.48
1:1:1350:C:C2'	1:1:1351:C:C6	2.82	0.48
1:1:1469:A:N7	1:1:1521:G:N2	2.60	0.48
1:1:1501:G:OP1	4:B:101:ARG:NH2	2.43	0.48
1:1:1682:G:OP2	1:1:1699:G:N2	2.39	0.48
1:1:2308:G:H2'	1:1:2310:C:H5	1.75	0.48
1:1:2391:G:H22	1:1:2424:C:H2'	1.77	0.48
2:2:254:G:O2'	47:v:18:GLU:O	2.26	0.48
2:2:1404:C:H42	2:2:1497:G:H1	1.60	0.48
40:o:66:GLU:OE2	40:o:68:ARG:NE	2.46	0.48
1:1:752:A:H62	1:1:2609:U:H3	1.61	0.48
1:1:1945:G:N2	1:1:1963:U:O4	2.46	0.48
1:1:2515:C:H2'	1:1:2516:A:H8	1.78	0.48
2:2:413:G:O3'	2:2:428:G:N2	2.46	0.48
2:2:1112:C:N4	33:h:176:HIS:O	2.46	0.48
2:2:1158:C:O2'	2:2:1160:G:O5'	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:77:U:OP1	22:V:21:ARG:NH1	2.46	0.48
24:X:36:HIS:ND1	24:X:37:ARG:O	2.46	0.48
32:g:120:GLN:HA	32:g:124:GLY:HA3	1.95	0.48
1:1:1755:A:H5'	16:P:103:ARG:HH22	1.78	0.48
2:2:243:A:H62	2:2:281:G:H1'	1.78	0.48
2:2:380:G:N2	2:2:383:A:OP2	2.32	0.48
2:2:440:C:H2'	2:2:441:A:H8	1.78	0.48
2:2:615:G:N1	2:2:626:G:C5	2.81	0.48
2:2:626:G:O3'	46:u:51:ARG:NH1	2.39	0.48
2:2:750:C:H4'	45:t:21:ASP:HA	1.96	0.48
2:2:1249:C:N4	2:2:1288:A:OP2	2.33	0.48
3:3:31:C:O2'	3:3:32:U:H5'	2.13	0.48
34:i:147:GLU:C	34:i:149:ALA:N	2.72	0.48
1:1:99:U:O2'	1:1:100:U:H5	1.94	0.48
1:1:572:A:OP2	18:R:80:ARG:NH2	2.46	0.48
1:1:1059:G:H3'	1:1:1060:U:H3'	1.96	0.48
1:1:1475:G:O2'	1:1:1514:G:O6	2.31	0.48
1:1:2047:C:H2'	1:1:2048:G:C8	2.48	0.48
9:G:100:ALA:O	9:G:104:THR:OG1	2.28	0.48
23:W:50:ASN:HB2	23:W:82:ILE:HB	1.95	0.48
35:j:38:VAL:HG21	35:j:114:VAL:HG23	1.95	0.48
48:w:42:SER:OG	48:w:47:THR:O	2.32	0.48
1:1:672:C:OP2	12:L:42:SER:OG	2.29	0.48
1:1:2394:C:H5''	12:L:63:LYS:HE2	1.95	0.48
2:2:845:A:H8	48:w:15:ALA:HB2	1.79	0.48
8:F:42:GLU:O	8:F:53:GLY:N	2.47	0.48
16:P:63:LYS:HE2	16:P:65:SER:HB2	1.95	0.48
1:1:558:U:H2'	1:1:559:G:H8	1.78	0.48
1:1:636:G:N1	12:L:76:GLU:OE1	2.38	0.48
1:1:1352:U:O2	1:1:1380:G:N2	2.46	0.48
1:1:1782:U:H1'	1:1:2609:U:H5''	1.96	0.48
1:1:1934:C:H2'	1:1:1935:G:H8	1.79	0.48
1:1:2163:A:OP2	1:1:2171:A:N6	2.40	0.48
2:2:437:U:O2'	34:i:120:HIS:ND1	2.37	0.48
2:2:824:G:H2'	2:2:825:A:H8	1.79	0.48
2:2:880:C:H2'	2:2:881:G:H8	1.79	0.48
2:2:1216:A:H5''	44:s:5:SER:HB3	1.95	0.48
38:m:39:VAL:O	38:m:43:GLU:N	2.42	0.48
47:v:47:HIS:N	47:v:73:TRP:O	2.45	0.48
1:1:717:C:H3'	1:1:718:A:H8	1.77	0.48
1:1:956:G:OP2	13:M:86:LYS:NZ	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1750:G:O2'	1:1:2860:A:N1	2.44	0.48
2:2:251:G:O6	2:2:272:C:N4	2.46	0.48
4:B:71:LYS:O	4:B:118:SER:OG	2.31	0.48
33:h:26:THR:HG23	44:s:76:LYS:HG3	1.95	0.48
33:h:79:LYS:HB2	33:h:82:GLU:HB3	1.96	0.48
36:k:36:ILE:HD11	36:k:39:LEU:HB2	1.95	0.48
1:1:910:A:N6	1:1:2277:G:O2'	2.47	0.48
1:1:1305:C:N4	1:1:1607:C:OP2	2.47	0.48
1:1:2653:U:O2	8:F:110:SER:OG	2.31	0.48
4:B:157:SER:OG	4:B:158:ALA:N	2.43	0.48
34:i:85:ASN:HB3	34:i:88:GLU:HB2	1.96	0.48
39:n:34:SER:HB3	39:n:37:GLN:HG3	1.94	0.48
44:s:47:LYS:O	44:s:50:THR:OG1	2.31	0.48
1:1:372:G:O2'	1:1:400:G:O6	2.30	0.48
1:1:510:C:H2'	1:1:511:U:O4'	2.14	0.48
1:1:995:C:O2	10:J:3:THR:OG1	2.30	0.48
1:1:1380:G:N3	1:1:1381:G:C8	2.82	0.48
1:1:2200:C:H42	1:1:2223:G:H1	1.60	0.48
1:1:2365:G:N7	30:e:39:LYS:NZ	2.44	0.48
1:1:2754:U:O3'	31:f:19:ARG:NH2	2.47	0.48
2:2:200:G:H2'	2:2:201:G:H8	1.79	0.48
2:2:675:A:O2'	41:p:116:ILE:O	2.31	0.48
2:2:694:A:OP1	41:p:55:SER:OG	2.27	0.48
2:2:845:A:H1'	48:w:11:CYS:HB3	1.95	0.48
9:G:14:SER:OG	9:G:15:LEU:N	2.47	0.48
32:g:20:THR:HA	32:g:39:HIS:HE1	1.78	0.48
33:h:110:GLU:HG2	33:h:140:ASN:HB3	1.95	0.48
54:A:248:HIS:HB2	54:A:274:LEU:HD21	1.96	0.48
1:1:52:A:OP2	1:1:117:G:N1	2.31	0.48
1:1:220:G:H22	1:1:427:U:H2'	1.78	0.48
1:1:647:G:H3'	1:1:648:G:C8	2.47	0.48
1:1:2064:C:H42	1:1:2446:G:H1	1.61	0.48
1:1:2626:C:H2'	1:1:2627:G:H8	1.79	0.48
2:2:269:C:H2'	2:2:270:A:C8	2.49	0.48
2:2:652:U:O4	2:2:752:G:O2'	2.28	0.48
2:2:837:U:H2'	2:2:838:G:H8	1.78	0.48
15:O:51:ALA:HB3	15:O:78:VAL:HB	1.96	0.48
27:b:39:LEU:HB2	27:b:42:HIS:HB2	1.95	0.48
35:j:96:MET:HA	35:j:125:ALA:HA	1.95	0.48
37:l:25:LYS:HA	37:l:28:ASN:HD22	1.79	0.48
37:l:64:VAL:O	37:l:68:ASN:ND2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:q:72:HIS:HB2	42:q:74:LEU:HD23	1.96	0.48
1:1:603:A:O4'	1:1:655:A:N6	2.47	0.47
1:1:2787:C:H1'	5:C:63:PRO:HG3	1.95	0.47
2:2:1055:A:O2'	33:h:156:ARG:NH2	2.47	0.47
2:2:1323:G:H2'	2:2:1324:A:C8	2.49	0.47
2:2:1391:U:H2'	2:2:1392:G:C8	2.49	0.47
21:U:86:ARG:NH1	21:U:100:SER:OG	2.47	0.47
34:i:108:GLY:HA3	34:i:114:ALA:HB2	1.96	0.47
44:s:3:LYS:HB2	44:s:6:MET:HG2	1.96	0.47
54:A:310:TYR:HE1	54:A:317:VAL:HG13	1.79	0.47
1:1:2320:U:O2	1:1:2333:A:N6	2.47	0.47
2:2:130:A:N1	2:2:233:C:O2'	2.46	0.47
2:2:519:C:N4	2:2:529:G:O2'	2.44	0.47
2:2:950:U:H3'	43:r:101:ARG:HH22	1.79	0.47
2:2:1438:G:H1	2:2:1463:U:H3	1.62	0.47
7:E:57:LEU:HD22	7:E:65:PRO:HB3	1.96	0.47
17:Q:82:GLY:O	17:Q:86:ALA:N	2.44	0.47
22:V:57:TYR:OH	22:V:79:ARG:NH2	2.47	0.47
1:1:938:G:H2'	1:1:939:G:H8	1.80	0.47
1:1:1468:U:H2'	1:1:1522:A:N6	2.29	0.47
1:1:2655:G:N2	1:1:2656:U:O4	2.37	0.47
33:h:112:ASP:HB3	33:h:115:LEU:HB2	1.96	0.47
36:k:1:MET:N	36:k:65:GLU:OE2	2.47	0.47
37:l:78:ARG:HG2	37:l:80:VAL:HG13	1.96	0.47
42:q:24:LEU:HG	42:q:30:LYS:HG2	1.96	0.47
1:1:538:A:H5''	10:J:7:LYS:HE2	1.96	0.47
1:1:1028:A:OP2	1:1:1126:A:N6	2.44	0.47
1:1:1073:A:H62	1:1:1074:G:H21	1.62	0.47
1:1:2830:C:O2'	1:1:2883:A:N1	2.40	0.47
5:C:151:THR:HG22	5:C:152:PRO:HD3	1.97	0.47
21:U:86:ARG:NH1	21:U:100:SER:O	2.36	0.47
28:c:14:SER:OG	28:c:48:ILE:O	2.29	0.47
34:i:56:ARG:HH12	34:i:63:ARG:HH12	1.63	0.47
40:o:64:GLN:O	44:s:99:ALA:N	2.46	0.47
44:s:48:LEU:HD12	44:s:51:LEU:HD12	1.97	0.47
1:1:1044:C:O2'	1:1:1111:A:N1	2.48	0.47
1:1:1604:C:O2'	1:1:1610:A:N1	2.47	0.47
1:1:2136:G:N2	1:1:2156:G:N3	2.63	0.47
2:2:6:G:H1	35:j:100:SER:H	1.62	0.47
2:2:543:U:OP1	34:i:14:ARG:NE	2.39	0.47
2:2:1403:C:N4	55:4:18:C:OP1	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:F:84:THR:OG1	8:F:134:LYS:NZ	2.40	0.47
40:o:32:THR:O	40:o:80:THR:OG1	2.32	0.47
41:p:15:GLN:NE2	41:p:78:GLY:O	2.47	0.47
54:A:153:GLU:CG	54:A:159:TYR:HE1	2.26	0.47
1:1:743:A:O2'	1:1:1659:G:OP1	2.28	0.47
1:1:1584:U:H4'	1:1:1585:C:H5''	1.95	0.47
1:1:2372:U:H2'	1:1:2373:G:C8	2.49	0.47
1:1:2522:U:O2'	1:1:2647:U:OP1	2.26	0.47
2:2:1097:C:O2'	2:2:1169:A:N3	2.41	0.47
2:2:1359:C:H3'	44:s:75:ARG:HH22	1.80	0.47
2:2:1441:A:H2'	2:2:1442:G:H8	1.79	0.47
21:U:8:ASP:N	21:U:8:ASP:OD1	2.46	0.47
34:i:60:LYS:HE3	34:i:195:ILE:HG12	1.96	0.47
39:n:35:LEU:HD11	39:n:45:ARG:HA	1.96	0.47
41:p:35:THR:OG1	41:p:36:ASP:N	2.47	0.47
50:y:67:ILE:HB	50:y:71:LYS:HD3	1.97	0.47
1:1:1047:G:N2	1:1:1110:G:O2'	2.48	0.47
1:1:1667:G:O2'	1:1:1991:U:O4	2.32	0.47
2:2:253:A:O2'	47:v:17:MET:SD	2.70	0.47
2:2:291:U:H2'	2:2:292:G:H8	1.79	0.47
2:2:880:C:H3'	42:q:6:GLN:HE21	1.78	0.47
2:2:977:A:N6	2:2:1224:U:O5'	2.45	0.47
2:2:1145:A:O2'	2:2:1146:A:O4'	2.33	0.47
2:2:1350:A:O2'	37:l:33:ASP:OD1	2.28	0.47
5:C:5:VAL:H	5:C:32:ASN:HD21	1.63	0.47
8:F:89:LEU:HD23	8:F:94:TYR:HB3	1.96	0.47
9:G:127:GLU:HA	9:G:145:ASN:HA	1.96	0.47
19:S:4:ILE:HG12	19:S:106:VAL:HG22	1.96	0.47
19:S:66:ILE:HA	19:S:69:LEU:HD13	1.96	0.47
32:g:15:HIS:HA	32:g:41:ILE:HB	1.95	0.47
32:g:183:VAL:H	32:g:197:ASP:HB2	1.79	0.47
33:h:70:THR:O	33:h:106:VAL:N	2.48	0.47
34:i:62:ARG:HH21	34:i:68:LEU:HA	1.80	0.47
41:p:82:LEU:O	41:p:108:THR:N	2.47	0.47
42:q:56:ARG:NE	42:q:62:GLU:OE1	2.37	0.47
48:w:33:ILE:HD12	48:w:37:GLY:HA2	1.96	0.47
49:x:3:ARG:NH1	49:x:9:PRO:O	2.48	0.47
51:z:5:LYS:O	51:z:18:ARG:NH2	2.47	0.47
52:5:15:C:H5'	52:5:16:C:C6	2.49	0.47
1:1:1006:C:O2'	10:J:108:MET:O	2.32	0.47
1:1:1919:A:N1	2:2:1495:U:O2'	2.42	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2620:C:O2'	5:C:162:ALA:O	2.32	0.47
2:2:405:U:O4	34:i:2:ALA:N	2.48	0.47
2:2:437:U:O2'	34:i:154:ARG:NH2	2.47	0.47
2:2:723:U:O4	51:z:56:HIS:ND1	2.47	0.47
2:2:1162:C:H2'	2:2:1163:A:H8	1.80	0.47
2:2:1442:G:H1	2:2:1461:G:H1'	1.79	0.47
4:B:145:GLU:OE1	4:B:149:GLY:N	2.45	0.47
16:P:54:GLY:N	16:P:57:SER:OG	2.48	0.47
17:Q:47:TYR:OH	17:Q:51:ARG:NH2	2.48	0.47
38:m:26:THR:HA	38:m:60:GLU:HA	1.96	0.47
40:o:10:LEU:HB2	40:o:72:ARG:HB2	1.96	0.47
41:p:60:PRO:HD3	41:p:91:PRO:HB2	1.96	0.47
47:v:12:VAL:HG22	47:v:23:VAL:HG22	1.97	0.47
47:v:26:GLU:HA	47:v:41:THR:HA	1.96	0.47
1:1:589:U:H2'	1:1:590:A:C8	2.50	0.47
1:1:1397:U:OP2	1:1:1398:C:N4	2.37	0.47
2:2:865:A:N3	2:2:918:A:O2'	2.45	0.47
2:2:1005:A:H3'	2:2:1006:G:C8	2.50	0.47
7:E:30:ARG:O	7:E:158:THR:OG1	2.30	0.47
34:i:69:GLU:O	34:i:73:ARG:N	2.44	0.47
50:y:35:VAL:HG21	50:y:54:MET:HG2	1.97	0.47
1:1:238:C:O2'	1:1:608:A:N3	2.44	0.47
2:2:1386:G:H2'	2:2:1387:G:H8	1.80	0.47
3:3:37:C:N3	3:3:48:U:O2'	2.41	0.47
16:P:25:THR:HB	16:P:88:ARG:HB3	1.97	0.47
39:n:55:VAL:HG11	39:n:94:LEU:HD13	1.97	0.47
54:A:108:GLU:HA	54:A:170:GLY:HA2	1.97	0.47
1:1:536:G:H1	1:1:557:C:H42	1.62	0.46
1:1:2039:U:H2'	1:1:2040:G:C8	2.50	0.46
1:1:2515:C:H2'	1:1:2516:A:C8	2.50	0.46
18:R:6:GLN:HG3	18:R:39:LEU:HD11	1.97	0.46
32:g:59:LYS:O	32:g:62:SER:OG	2.30	0.46
37:l:146:GLU:HA	37:l:149:LYS:HB2	1.95	0.46
41:p:31:ILE:HG12	41:p:46:THR:HG22	1.95	0.46
54:A:180:GLY:H	54:A:203:VAL:HB	1.81	0.46
1:1:251:A:OP1	12:L:58:TYR:OH	2.28	0.46
1:1:1209:U:O2'	1:1:1237:A:N1	2.43	0.46
1:1:1338:G:H2'	1:1:1339:G:C8	2.50	0.46
1:1:1595:C:H2'	1:1:1596:A:H8	1.79	0.46
1:1:1721:G:H22	1:1:1738:G:H2'	1.79	0.46
1:1:2000:C:O2'	1:1:2001:C:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2072:C:H42	1:1:2437:G:H1	1.63	0.46
2:2:1257:A:O2'	2:2:1258:G:H5'	2.15	0.46
2:2:1320:C:O2	49:x:36:ARG:NH2	2.48	0.46
12:L:73:ILE:HD12	12:L:106:GLU:HB2	1.96	0.46
1:1:926:G:H2'	1:1:927:A:H8	1.81	0.46
1:1:1106:G:H2'	1:1:1107:G:H8	1.79	0.46
2:2:390:U:O2'	46:u:28:ARG:NH2	2.44	0.46
2:2:494:G:H2'	2:2:496:A:H8	1.80	0.46
2:2:886:G:H1	2:2:911:U:H3	1.60	0.46
2:2:1395:C:HO2'	2:2:1401:G:HO2'	1.60	0.46
11:K:43:ILE:HD12	11:K:56:ASP:HB2	1.96	0.46
47:v:10:GLY:N	47:v:59:VAL:O	2.44	0.46
1:1:377:G:H1	1:1:397:U:H3	1.64	0.46
1:1:715:A:H2	45:t:40:GLN:HE22	1.62	0.46
1:1:729:G:C5	4:B:207:LYS:HB2	2.50	0.46
2:2:855:U:OP2	2:2:871:U:N3	2.38	0.46
2:2:1013:G:N2	2:2:1016:A:OP2	2.43	0.46
14:N:28:LEU:O	14:N:32:GLU:N	2.39	0.46
32:g:137:ARG:O	32:g:141:LEU:N	2.43	0.46
54:A:139:ALA:HB1	54:A:144:TRP:HB2	1.98	0.46
54:A:177:PHE:HB3	54:A:321:ARG:HH22	1.80	0.46
1:1:1041:G:N2	1:1:1042:G:O6	2.48	0.46
1:1:1967:C:H2'	1:1:1968:G:O4'	2.16	0.46
2:2:309:A:H2'	2:2:310:G:H8	1.80	0.46
2:2:1072:G:O6	2:2:1102:A:N6	2.48	0.46
2:2:1438:G:N2	2:2:1463:U:O2	2.39	0.46
2:2:1512:U:H2'	2:2:1513:A:C8	2.51	0.46
4:B:93:LEU:HD21	4:B:101:ARG:HD3	1.98	0.46
6:D:31:VAL:HG21	6:D:104:ALA:HB2	1.97	0.46
35:j:102:GLY:H	35:j:122:ASN:ND2	2.14	0.46
37:l:141:VAL:O	37:l:145:ALA:N	2.48	0.46
49:x:15:LEU:HD13	49:x:33:THR:HG21	1.96	0.46
1:1:692:C:HO2'	1:1:1354:A:HO2'	1.57	0.46
1:1:1314:C:H42	1:1:1338:G:H1	1.63	0.46
1:1:2314:A:O2'	7:E:157:THR:OG1	2.34	0.46
2:2:177:G:OP2	50:y:64:LYS:NZ	2.48	0.46
7:E:48:LYS:HA	7:E:51:ASP:HB3	1.97	0.46
15:O:30:ARG:HD3	15:O:102:ARG:HH12	1.81	0.46
34:i:124:MET:HG3	34:i:146:ARG:HB3	1.96	0.46
34:i:172:GLU:HB2	34:i:183:LYS:HD2	1.97	0.46
36:k:88:MET:HE2	36:k:90:MET:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:l:138:ARG:HE	37:l:139:GLU:HG2	1.81	0.46
1:1:99:U:H5''	1:1:100:U:H5'	1.98	0.46
1:1:2233:U:H2'	1:1:2234:G:C8	2.51	0.46
1:1:2753:A:O2'	31:f:15:LYS:NZ	2.33	0.46
2:2:1004:A:O2'	2:2:1036:A:N6	2.43	0.46
4:B:62:TYR:HA	4:B:86:ASN:HD21	1.81	0.46
6:D:83:VAL:HB	6:D:86:ALA:HB2	1.96	0.46
8:F:95:ARG:HB2	8:F:128:GLN:HB3	1.98	0.46
28:c:10:LYS:HA	28:c:22:THR:HA	1.98	0.46
40:o:26:VAL:HG13	40:o:36:VAL:HG11	1.96	0.46
54:A:330:ASP:N	54:A:330:ASP:OD1	2.47	0.46
1:1:1453:A:O2'	14:N:63:ARG:NH1	2.48	0.46
1:1:1693:U:O4	1:1:1976:U:O2'	2.31	0.46
1:1:2036:C:H2'	1:1:2037:A:C8	2.51	0.46
1:1:2375:G:N2	1:1:2378:A:OP2	2.38	0.46
1:1:2892:G:H5''	1:1:2893:A:H5'	1.98	0.46
2:2:417:G:H1	2:2:426:U:H3	1.63	0.46
2:2:501:C:OP2	42:q:121:ARG:NH1	2.49	0.46
34:i:187:GLU:H	34:i:190:ASP:HB2	1.80	0.46
37:l:113:ASP:OD2	37:l:122:ASN:ND2	2.41	0.46
52:5:11:G:N2	52:5:24:U:O2	2.49	0.46
1:1:265:A:O2'	1:1:428:A:N6	2.49	0.46
1:1:2566:A:N1	11:K:28:SER:OG	2.46	0.46
2:2:294:U:OP1	2:2:610:U:O2'	2.27	0.46
2:2:407:U:OP1	34:i:3:ARG:NH2	2.36	0.46
2:2:868:C:H2'	2:2:869:G:O4'	2.16	0.46
2:2:1304:G:N2	2:2:1332:A:OP2	2.37	0.46
17:Q:17:ILE:HG13	17:Q:32:TYR:HE1	1.80	0.46
30:e:56:GLY:HA2	30:e:59:ILE:HD12	1.98	0.46
33:h:71:ALA:HA	33:h:106:VAL:HB	1.98	0.46
43:r:75:MET:O	43:r:79:ARG:N	2.41	0.46
1:1:309:A:N3	1:1:329:G:O2'	2.47	0.46
35:j:111:MET:HA	35:j:114:VAL:HG12	1.97	0.46
38:m:92:LEU:O	38:m:117:ARG:NH2	2.38	0.46
40:o:28:THR:O	40:o:32:THR:OG1	2.33	0.46
45:t:21:ASP:OD1	45:t:21:ASP:N	2.49	0.46
1:1:613:A:C2	1:1:614:A:C4	3.04	0.45
1:1:2241:A:H2'	1:1:2242:G:C8	2.51	0.45
2:2:2:A:N3	2:2:613:C:O2'	2.38	0.45
2:2:257:G:H2'	2:2:258:G:C8	2.52	0.45
4:B:108:LYS:HA	4:B:196:GLY:HA2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:e:33:LEU:HB3	30:e:41:LYS:HE2	1.97	0.45
43:r:2:ALA:O	43:r:9:ILE:N	2.44	0.45
1:1:57:C:H2'	1:1:58:G:O4'	2.16	0.45
1:1:340:A:O2'	6:D:162:ARG:NH1	2.50	0.45
1:1:861:A:N3	3:3:79:G:O2'	2.48	0.45
1:1:1381:G:O2'	1:1:1572:A:N1	2.49	0.45
1:1:1982:U:H2'	1:1:1983:G:H8	1.80	0.45
1:1:2313:C:O4'	7:E:37:ASN:ND2	2.49	0.45
1:1:2477:U:O2	31:f:4:ARG:NH2	2.38	0.45
1:1:2808:G:O2'	1:1:2890:G:O6	2.29	0.45
2:2:1:A:OP2	2:2:628:G:O2'	2.28	0.45
2:2:1071:C:H2'	2:2:1072:G:H8	1.80	0.45
2:2:1124:G:O2'	2:2:1127:G:N1	2.49	0.45
2:2:1471:U:H2'	2:2:1472:U:C6	2.51	0.45
6:D:1:MET:HB3	6:D:14:VAL:HG23	1.98	0.45
7:E:117:LEU:N	7:E:177:PHE:O	2.50	0.45
33:h:58:GLU:O	33:h:65:ARG:N	2.45	0.45
43:r:20:THR:HA	43:r:25:VAL:HG23	1.99	0.45
47:v:59:VAL:HG23	47:v:78:VAL:HG22	1.99	0.45
51:z:43:THR:HA	51:z:46:LYS:HB3	1.98	0.45
1:1:407:G:H2'	1:1:408:G:H8	1.81	0.45
1:1:2526:G:H1	1:1:2537:U:H3	1.64	0.45
2:2:396:C:O2'	2:2:398:U:OP1	2.34	0.45
2:2:542:G:H5'	34:i:39:GLY:HA3	1.97	0.45
2:2:651:C:N4	2:2:753:A:OP2	2.44	0.45
2:2:948:C:H2'	2:2:949:A:H8	1.81	0.45
2:2:1359:C:O5'	44:s:75:ARG:NH2	2.50	0.45
4:B:171:TYR:HE1	4:B:185:GLU:HG2	1.80	0.45
8:F:86:LYS:HG2	8:F:132:VAL:HG22	1.97	0.45
15:O:53:THR:HB	15:O:65:THR:HB	1.98	0.45
17:Q:26:GLY:O	17:Q:30:ARG:NH1	2.49	0.45
26:Z:4:THR:HA	26:Z:39:GLU:HA	1.97	0.45
1:1:611:C:H2'	1:1:612:G:O4'	2.16	0.45
1:1:1564:C:OP2	4:B:26:LYS:NZ	2.41	0.45
1:1:1845:G:H2'	1:1:1846:G:C8	2.52	0.45
1:1:2576:G:O2'	1:1:2579:C:OP2	2.28	0.45
1:1:2676:C:O2	1:1:2732:G:N2	2.49	0.45
2:2:41:G:H2'	2:2:42:G:H8	1.81	0.45
2:2:123:U:OP1	2:2:311:C:O2'	2.33	0.45
2:2:1166:G:C2	2:2:1168:U:H4'	2.51	0.45
17:Q:50:ARG:O	17:Q:54:LYS:NZ	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:c:53:LYS:HG2	28:c:55:LYS:H	1.81	0.45
46:u:4:ILE:HG12	46:u:21:VAL:HG22	1.97	0.45
1:1:143:C:H3'	1:1:144:A:H8	1.82	0.45
1:1:188:G:H5'	24:X:14:THR:HG21	1.99	0.45
1:1:226:A:O2'	1:1:229:C:N4	2.50	0.45
1:1:321:U:H5''	6:D:131:THR:HG23	1.98	0.45
1:1:1170:C:H2'	1:1:1171:G:H8	1.82	0.45
2:2:501:C:H2'	2:2:502:A:C8	2.52	0.45
2:2:770:C:O2'	2:2:899:C:N3	2.46	0.45
2:2:877:G:H2'	2:2:878:A:H8	1.81	0.45
16:P:53:ARG:N	16:P:57:SER:OG	2.45	0.45
39:n:8:GLY:N	39:n:19:VAL:O	2.44	0.45
43:r:68:ASP:OD1	43:r:71:ARG:NH1	2.38	0.45
51:z:41:PRO:HA	51:z:44:GLU:HB3	1.98	0.45
52:5:58:A:O2'	52:5:60:U:OP2	2.25	0.45
1:1:603:A:O5'	1:1:655:A:N6	2.50	0.45
2:2:187:G:C2	2:2:189:A:H5''	2.51	0.45
2:2:615:G:C5	2:2:626:G:O6	2.70	0.45
2:2:1376:U:O4	37:l:10:ARG:NH1	2.39	0.45
15:O:52:SER:OG	15:O:53:THR:N	2.50	0.45
16:P:71:GLU:OE2	16:P:101:ARG:NE	2.50	0.45
39:n:12:ARG:HG3	39:n:13:LYS:H	1.82	0.45
1:1:414:C:H2'	1:1:415:A:C8	2.52	0.45
1:1:729:G:N3	1:1:729:G:C2'	2.73	0.45
1:1:874:G:H2'	1:1:875:G:H8	1.82	0.45
1:1:2627:G:N2	1:1:2777:G:OP2	2.50	0.45
2:2:299:G:N2	2:2:565:U:O2	2.49	0.45
3:3:31:C:H1'	3:3:53:A:C6	2.52	0.45
7:E:158:THR:OG1	7:E:159:THR:N	2.48	0.45
26:Z:9:GLN:HB2	26:Z:29:LEU:HD13	1.99	0.45
33:h:123:GLN:HB3	33:h:128:VAL:HG21	1.99	0.45
41:p:36:ASP:OD1	41:p:40:ASN:N	2.50	0.45
43:r:16:VAL:HA	43:r:19:LEU:HB2	1.99	0.45
1:1:210:C:O2'	1:1:1366:A:O2'	2.34	0.45
1:1:1386:C:H2'	1:1:1387:A:C8	2.52	0.45
1:1:2114:A:N6	1:1:2167:U:O2'	2.50	0.45
14:N:55:ALA:HA	14:N:80:PHE:CE1	2.51	0.45
28:c:17:THR:HG22	28:c:19:HIS:H	1.81	0.45
33:h:40:ARG:NH1	33:h:55:ILE:O	2.40	0.45
39:n:120:LYS:HB2	39:n:123:ARG:HB3	1.99	0.45
45:t:26:GLU:HG3	45:t:81:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:x:50:ALA:HA	49:x:59:PRO:HA	1.98	0.45
1:1:692:C:O2'	1:1:1354:A:O2'	2.29	0.45
1:1:2010:G:H5''	19:S:42:LYS:HB2	1.98	0.45
1:1:2828:G:H2'	1:1:2829:A:H8	1.81	0.45
2:2:21:G:H2'	2:2:22:G:C8	2.52	0.45
2:2:461:A:H2'	2:2:462:G:C8	2.51	0.45
2:2:522:C:OP2	42:q:66:TYR:OH	2.31	0.45
2:2:715:A:H2'	2:2:716:A:C8	2.52	0.45
2:2:755:G:H21	38:m:4:GLN:HE22	1.64	0.45
2:2:946:A:H2'	2:2:947:G:C8	2.52	0.45
28:c:8:LYS:HA	28:c:24:THR:HA	1.98	0.45
38:m:18:GLN:HG3	38:m:72:VAL:HB	1.99	0.45
1:1:729:G:OP1	4:B:10:SER:OG	2.35	0.45
1:1:2189:U:H1'	1:1:2190:G:N7	2.32	0.45
2:2:1209:C:O2'	2:2:1214:C:N4	2.40	0.45
33:h:114:LYS:HB2	33:h:185:ASN:HD22	1.82	0.45
1:1:1380:G:O6	1:1:1381:G:C6	2.66	0.44
1:1:1538:G:H2'	1:1:1539:U:C6	2.52	0.44
2:2:321:A:H3'	2:2:328:C:H41	1.82	0.44
2:2:1308:U:H4'	43:r:91:HIS:HE1	1.81	0.44
3:3:31:C:H2'	3:3:32:U:C6	2.52	0.44
1:1:2667:C:N3	8:F:110:SER:OG	2.50	0.44
2:2:41:G:H2'	2:2:42:G:C8	2.53	0.44
2:2:1412:C:H2'	2:2:1413:A:C8	2.51	0.44
2:2:1442:G:N2	2:2:1461:G:O2'	2.50	0.44
3:3:39:A:O2'	3:3:46:A:N1	2.48	0.44
5:C:25:THR:OG1	5:C:191:GLY:O	2.30	0.44
21:U:11:VAL:HA	21:U:73:PHE:H	1.82	0.44
34:i:98:LEU:HB2	34:i:135:TYR:HB3	1.99	0.44
43:r:50:GLU:O	43:r:54:ASP:N	2.44	0.44
44:s:5:SER:O	44:s:9:ARG:N	2.44	0.44
53:H:78:PHE:O	53:H:78:PHE:CD1	2.70	0.44
1:1:33:C:O2	1:1:447:A:N6	2.49	0.44
1:1:510:C:H5'	1:1:510:C:H6	1.81	0.44
1:1:1170:C:N4	1:1:1171:G:O6	2.51	0.44
1:1:1350:C:C2	1:1:1351:C:C5	3.06	0.44
1:1:2037:A:H2'	1:1:2038:G:C8	2.53	0.44
2:2:231:U:H2'	2:2:232:G:C8	2.53	0.44
2:2:620:C:H2'	2:2:621:A:O4'	2.17	0.44
2:2:1020:G:H3'	2:2:1021:A:H8	1.81	0.44
4:B:107:PRO:HD2	4:B:110:LEU:HD22	1.97	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:U:11:VAL:HG12	21:U:72:ILE:HA	1.99	0.44
33:h:155:GLY:HA2	33:h:163:ALA:HB1	1.99	0.44
37:l:72:THR:OG1	37:l:142:HIS:NE2	2.39	0.44
1:1:244:A:H62	1:1:254:G:H21	1.65	0.44
1:1:1041:G:N2	1:1:1114:C:N3	2.66	0.44
1:1:2390:U:O5'	30:e:35:LYS:NZ	2.45	0.44
2:2:103:U:O2'	2:2:171:A:N1	2.45	0.44
2:2:408:A:OP1	34:i:110:THR:OG1	2.33	0.44
2:2:1329:A:H5''	43:r:25:VAL:HA	1.99	0.44
4:B:37:ASN:HB2	4:B:62:TYR:HB2	1.99	0.44
39:n:23:PRO:HA	39:n:61:LEU:HA	1.98	0.44
1:1:598:U:H2'	1:1:599:A:C8	2.52	0.44
1:1:989:G:H5''	26:Z:14:ILE:HD11	1.99	0.44
1:1:1350:C:N3	1:1:1382:G:C2	2.85	0.44
1:1:1859:U:H2'	1:1:1860:G:O4'	2.18	0.44
1:1:2339:C:H5''	1:1:2339:C:C6	2.52	0.44
1:1:2627:G:O2'	1:1:2781:A:N1	2.45	0.44
2:2:443:C:H2'	2:2:444:G:C8	2.53	0.44
2:2:460:A:H3'	2:2:461:A:H8	1.83	0.44
2:2:672:U:O2'	48:w:64:TYR:OH	2.33	0.44
2:2:959:A:HO2'	2:2:984:C:HO2'	1.61	0.44
2:2:1359:C:OP2	44:s:75:ARG:NE	2.50	0.44
3:3:63:C:H2'	3:3:64:G:C8	2.52	0.44
18:R:52:PRO:HB2	18:R:53:PHE:CD1	2.52	0.44
33:h:70:THR:HG21	33:h:76:VAL:HG21	2.00	0.44
34:i:62:ARG:O	34:i:66:GLY:N	2.44	0.44
47:v:9:GLN:HA	47:v:60:GLU:HA	1.99	0.44
1:1:117:G:OP2	1:1:119:A:O2'	2.33	0.44
1:1:1073:A:C5	1:1:1074:G:H1'	2.53	0.44
1:1:1798:U:OP2	4:B:271:ARG:NH2	2.46	0.44
2:2:1158:C:O2'	2:2:1160:G:P	2.76	0.44
2:2:1158:C:O2	2:2:1158:C:H2'	2.17	0.44
2:2:1220:G:P	44:s:53:ARG:HH12	2.41	0.44
3:3:22:U:H2'	3:3:23:G:C8	2.52	0.44
23:W:37:ILE:HD11	23:W:61:ALA:HB2	1.99	0.44
32:g:32:PHE:HB2	32:g:42:ASN:HA	2.00	0.44
37:l:47:LEU:HD23	37:l:58:GLU:HB3	1.98	0.44
43:r:101:ARG:HG3	43:r:104:THR:H	1.81	0.44
50:y:4:ILE:HG22	50:y:6:SER:H	1.82	0.44
1:1:748:G:C8	1:1:750:A:C8	3.05	0.44
1:1:874:G:H2'	1:1:875:G:C8	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:882:G:N2	1:1:897:C:OP2	2.50	0.44
1:1:2059:A:C8	1:1:2503:2MA:HM22	2.52	0.44
1:1:2821:A:O3'	5:C:167:ASN:ND2	2.51	0.44
2:2:79:G:N2	2:2:90:C:O2	2.51	0.44
2:2:599:C:H2'	2:2:600:A:H8	1.82	0.44
2:2:615:G:C6	2:2:626:G:O6	2.71	0.44
2:2:751:U:H4'	45:t:24:SER:HA	1.99	0.44
2:2:1266:G:N2	2:2:1269:A:OP2	2.35	0.44
14:N:100:CYS:HA	27:b:43:ILE:HG12	2.00	0.44
19:S:20:VAL:HG21	19:S:43:ALA:HB3	2.00	0.44
19:S:29:VAL:HB	19:S:55:ILE:HD11	2.00	0.44
40:o:40:ILE:HB	40:o:73:LEU:HB3	1.99	0.44
54:A:134:MET:O	54:A:138:TYR:N	2.50	0.44
54:A:248:HIS:CE1	54:A:250:PRO:HD2	2.53	0.44
1:1:654:A:C4'	1:1:655:A:OP1	2.66	0.44
1:1:705:A:OP2	1:1:725:G:N2	2.49	0.44
1:1:1405:U:H2'	1:1:1406:U:O4'	2.18	0.44
1:1:1434:A:H2'	1:1:1435:G:H8	1.82	0.44
1:1:1449:G:H2'	1:1:1450:G:C8	2.53	0.44
1:1:1860:G:N2	1:1:1882:U:O2	2.39	0.44
1:1:2521:C:H2'	1:1:2522:U:C6	2.53	0.44
2:2:452:A:H62	2:2:480:U:H3	1.65	0.44
5:C:3:GLY:HA3	5:C:204:LYS:HG2	2.00	0.44
23:W:33:ALA:N	23:W:64:ASP:OD1	2.50	0.44
24:X:2:SER:HG	24:X:3:ARG:H	1.63	0.44
32:g:7:ARG:O	32:g:11:LYS:N	2.49	0.44
34:i:91:LEU:HA	34:i:94:LEU:HB2	1.99	0.44
36:k:33:GLU:O	36:k:35:LYS:NZ	2.50	0.44
37:l:77:SER:OG	37:l:86:GLN:OE1	2.29	0.44
54:A:223:ARG:N	54:A:247:THR:O	2.46	0.44
1:1:300:A:H1'	1:1:319:G:H1'	2.00	0.44
1:1:587:C:OP1	12:L:21:ARG:NH1	2.47	0.44
1:1:1754:A:O3'	16:P:103:ARG:NH2	2.51	0.44
1:1:1796:U:H3	1:1:1823:G:H1	1.66	0.44
1:1:2237:G:O2'	1:1:2239:G:N7	2.39	0.44
1:1:2831:G:O2'	1:1:2884:U:OP1	2.29	0.44
2:2:591:U:OP2	38:m:31:LYS:NZ	2.45	0.44
2:2:664:G:H22	2:2:741:G:H22	1.64	0.44
2:2:963:G:H2'	2:2:964:A:H8	1.83	0.44
2:2:1239:A:O2'	2:2:1298:U:O4	2.31	0.44
9:G:76:GLU:HG3	9:G:143:ILE:H	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:G:128:HIS:O	9:G:144:VAL:N	2.51	0.44
13:M:70:ASP:OD1	13:M:70:ASP:N	2.51	0.44
17:Q:39:VAL:O	17:Q:43:GLY:N	2.47	0.44
37:l:68:ASN:O	37:l:138:ARG:NH1	2.51	0.44
45:t:82:ILE:HG22	45:t:87:LEU:HD12	1.99	0.44
54:A:116:ARG:HA	54:A:161:GLU:HA	2.00	0.44
1:1:224:U:O4	1:1:419:U:O2'	2.36	0.43
1:1:323:C:H2'	6:D:163:ASN:HD21	1.81	0.43
1:1:635:C:O2'	1:1:639:U:OP1	2.31	0.43
1:1:1912:A:N1	2:2:1407:5MC:O2'	2.45	0.43
1:1:2315:G:OP1	7:E:33:LYS:NZ	2.47	0.43
2:2:858:G:H1	2:2:869:G:H3'	1.82	0.43
2:2:1048:G:H5''	44:s:3:LYS:HG2	2.00	0.43
2:2:1183:U:O2'	2:2:1185:G:OP2	2.35	0.43
30:e:28:ASN:O	30:e:36:LYS:NZ	2.51	0.43
38:m:52:GLU:O	38:m:58:GLU:N	2.46	0.43
44:s:46:LEU:HD12	49:x:13:LEU:HD13	2.00	0.43
1:1:687:C:H5''	29:d:2:LYS:HE2	2.00	0.43
1:1:1490:A:H3'	1:1:1491:G:H8	1.83	0.43
1:1:2549:G:H2'	1:1:2550:G:H8	1.83	0.43
1:1:2575:C:H5'	5:C:149:ASN:HB2	2.00	0.43
1:1:2699:C:H2'	1:1:2700:A:C8	2.53	0.43
2:2:335:C:H2'	2:2:336:A:H8	1.82	0.43
2:2:689:C:H2'	2:2:690:G:O4'	2.18	0.43
2:2:1069:C:H42	2:2:1106:G:H1	1.65	0.43
2:2:1239:A:O2'	37:l:114:LYS:O	2.36	0.43
9:G:89:LYS:HD3	9:G:123:ARG:HH22	1.83	0.43
26:Z:24:LEU:HD11	26:Z:54:MET:HE2	2.00	0.43
34:i:86:THR:HA	34:i:89:ASN:HB2	2.00	0.43
37:l:138:ARG:NE	37:l:139:GLU:OE2	2.51	0.43
1:1:26:G:N1	1:1:513:A:OP2	2.47	0.43
1:1:282:A:N3	1:1:359:G:N2	2.66	0.43
1:1:318:C:H2'	1:1:319:G:C8	2.54	0.43
1:1:741:U:H2'	1:1:742:A:C8	2.53	0.43
1:1:843:G:H2'	1:1:844:A:H8	1.84	0.43
1:1:1527:G:H21	1:1:1545:A:H62	1.66	0.43
1:1:2130:U:H4'	1:1:2158:A:H61	1.84	0.43
2:2:3:A:N1	2:2:629:A:H1'	2.33	0.43
8:F:101:ASN:HB2	8:F:117:LEU:HB2	2.00	0.43
8:F:117:LEU:HD12	8:F:121:ILE:HG21	2.00	0.43
39:n:51:PRO:HB3	39:n:84:THR:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:q:34:CYS:HB2	42:q:78:SER:H	1.82	0.43
1:1:612:G:H2'	1:1:614:A:C8	2.54	0.43
1:1:1779:U:OP2	1:1:1784:A:N6	2.52	0.43
1:1:2126:A:N6	1:1:2163:A:OP1	2.36	0.43
2:2:376:G:H5''	46:u:5:ARG:HB2	2.00	0.43
2:2:401:C:O2'	2:2:621:A:N3	2.49	0.43
2:2:980:C:O2'	44:s:13:ARG:NH1	2.52	0.43
2:2:1005:A:N6	2:2:1024:G:O2'	2.52	0.43
24:X:33:LEU:HA	24:X:52:SER:HA	1.99	0.43
52:5:65:C:O2'	52:5:66:C:H5'	2.19	0.43
1:1:5:A:H2'	1:1:6:A:C8	2.53	0.43
1:1:175:G:H2'	1:1:176:A:C8	2.53	0.43
1:1:658:U:O2'	6:D:97:ASN:OD1	2.36	0.43
1:1:1364:G:N2	1:1:1367:A:OP2	2.46	0.43
1:1:1595:C:H2'	1:1:1596:A:C8	2.54	0.43
1:1:1837:C:O2'	1:1:1927:A:N3	2.45	0.43
2:2:436:C:H4'	34:i:153:SER:HB3	1.99	0.43
2:2:696:A:N3	2:2:786:G:O2'	2.40	0.43
2:2:1342:C:H2'	2:2:1343:G:C8	2.54	0.43
18:R:68:ARG:O	18:R:90:ARG:NH2	2.50	0.43
32:g:24:ASN:HD22	32:g:192:ASP:HA	1.84	0.43
37:l:146:GLU:HG3	37:l:149:LYS:HD2	2.00	0.43
1:1:893:C:C2'	1:1:894:U:C6	3.02	0.43
1:1:1398:C:H2'	1:1:1399:C:C6	2.54	0.43
1:1:1413:A:N1	1:1:1589:U:N3	2.67	0.43
2:2:824:G:H2'	2:2:825:A:C8	2.54	0.43
2:2:1006:G:H22	2:2:1023:U:H2'	1.83	0.43
3:3:42:C:O2'	7:E:63:GLN:HB3	2.18	0.43
8:F:118:PRO:HD2	8:F:121:ILE:HG13	2.01	0.43
15:O:82:ALA:O	15:O:86:GLY:N	2.52	0.43
24:X:8:THR:OG1	24:X:9:GLY:N	2.51	0.43
44:s:40:ASP:O	44:s:44:ALA:N	2.51	0.43
1:1:759:G:H2'	1:1:760:G:H8	1.83	0.43
1:1:1166:G:N2	1:1:1183:U:O2	2.44	0.43
1:1:1246:A:H4'	6:D:40:ARG:HH12	1.82	0.43
1:1:2141:G:H2'	1:1:2142:A:C8	2.54	0.43
1:1:2243:U:H2'	1:1:2244:U:H6	1.84	0.43
1:1:2590:A:O3'	4:B:238:ARG:NH1	2.52	0.43
1:1:2756:U:OP2	31:f:19:ARG:NE	2.35	0.43
2:2:231:U:H2'	2:2:232:G:H8	1.84	0.43
2:2:362:G:O2'	2:2:364:A:N7	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:765:G:N2	2:2:813:U:OP2	2.52	0.43
2:2:844:G:C2	2:2:846:G:H1'	2.54	0.43
2:2:1136:C:H2'	2:2:1138:G:C6	2.54	0.43
3:3:79:G:H5''	13:M:18:ARG:HH22	1.83	0.43
5:C:109:VAL:HG22	5:C:203:VAL:HG22	2.00	0.43
35:j:18:VAL:HG21	35:j:56:VAL:HG13	2.00	0.43
1:1:183:C:O2'	1:1:432:A:N3	2.51	0.43
1:1:300:A:O5'	21:U:82:ARG:NH1	2.46	0.43
1:1:394:C:N4	1:1:395:U:O4	2.52	0.43
1:1:467:G:OP2	29:d:34:ARG:NH1	2.44	0.43
1:1:784:G:C2	4:B:228:VAL:HG11	2.54	0.43
1:1:1231:U:H2'	1:1:1232:G:H8	1.83	0.43
1:1:1548:A:H2'	1:1:1549:A:C8	2.54	0.43
1:1:1643:G:N2	1:1:1644:C:O2	2.51	0.43
1:1:1795:C:H2'	1:1:1796:U:O4'	2.19	0.43
1:1:1799:G:C5	4:B:176:LEU:HD13	2.54	0.43
1:1:2377:A:O2'	15:O:117:PHE:O	2.30	0.43
2:2:1118:U:H1'	2:2:1179:A:C4	2.53	0.43
1:1:845:A:N7	1:1:932:U:N3	2.58	0.43
1:1:932:U:H5'	1:1:933:A:C8	2.53	0.43
1:1:1026:G:H2'	1:1:1027:A:C8	2.52	0.43
1:1:1047:G:O2'	1:1:1110:G:N1	2.48	0.43
1:1:2521:C:O2'	1:1:2564:A:N3	2.47	0.43
2:2:565:U:H3'	2:2:566:G:H2'	2.01	0.43
2:2:1152:A:H5'	40:o:15:HIS:CD2	2.53	0.43
2:2:1269:A:N1	2:2:1312:G:O2'	2.49	0.43
2:2:1405:G:H2'	2:2:1406:U:H5''	2.01	0.43
8:F:166:ASP:OD1	8:F:166:ASP:N	2.51	0.43
9:G:47:PHE:O	9:G:51:ARG:N	2.52	0.43
11:K:34:GLY:N	11:K:37:ASP:OD2	2.52	0.43
16:P:88:ARG:HH12	16:P:110:ILE:HG13	1.83	0.43
27:b:31:ASP:OD2	27:b:34:SER:N	2.39	0.43
28:c:37:LYS:HG2	28:c:46:HIS:HB3	2.00	0.43
38:m:112:THR:HG23	38:m:115:ALA:H	1.84	0.43
39:n:84:THR:HG21	39:n:103:PHE:HB3	1.99	0.43
41:p:37:ARG:NH2	41:p:83:GLU:OE2	2.51	0.43
44:s:36:ALA:HB3	44:s:41:ARG:HD2	2.00	0.43
1:1:1350:C:H2'	1:1:1351:C:C5	2.48	0.43
1:1:1438:U:H2'	1:1:1439:A:H8	1.83	0.43
1:1:2372:U:H2'	1:1:2373:G:H8	1.84	0.43
1:1:2885:G:N7	27:b:40:ARG:NH1	2.55	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:214:C:H2'	2:2:215:C:H6	1.84	0.43
2:2:337:G:H2'	2:2:338:A:C8	2.54	0.43
2:2:832:G:N2	2:2:854:U:O2	2.41	0.43
2:2:1120:C:H2'	2:2:1121:U:C6	2.54	0.43
2:2:1280:A:OP2	40:o:9:ARG:NH2	2.52	0.43
34:i:57:GLU:HG3	34:i:200:ILE:HG13	2.01	0.43
42:q:50:ARG:HG3	42:q:90:LEU:HD21	2.00	0.43
45:t:5:THR:O	45:t:8:THR:OG1	2.28	0.43
54:A:144:TRP:CE2	54:A:171:VAL:HG22	2.54	0.43
1:1:107:G:H2'	1:1:108:G:H8	1.84	0.42
2:2:150:U:H2'	2:2:151:A:H8	1.84	0.42
2:2:403:C:H2'	2:2:404:G:C8	2.52	0.42
2:2:750:C:H2'	2:2:751:U:C6	2.53	0.42
2:2:965:U:H1'	2:2:969:A:C4	2.54	0.42
2:2:1320:C:H2'	2:2:1321:U:O4'	2.18	0.42
4:B:176:LEU:HD23	4:B:176:LEU:HA	1.89	0.42
6:D:23:PHE:HE1	6:D:108:ILE:HG22	1.84	0.42
12:L:2:ARG:O	12:L:5:THR:OG1	2.32	0.42
24:X:59:ILE:HG12	24:X:67:VAL:HG21	2.01	0.42
33:h:138:VAL:HA	33:h:149:ILE:HD13	2.01	0.42
36:k:62:MET:HG2	36:k:64:VAL:HG23	2.00	0.42
42:q:74:LEU:HD13	42:q:74:LEU:HA	1.91	0.42
43:r:67:GLY:O	43:r:71:ARG:N	2.50	0.42
1:1:154:U:H2'	1:1:155:A:C8	2.54	0.42
1:1:237:C:O2'	1:1:609:A:O2'	2.29	0.42
1:1:2885:G:O6	27:b:29:SER:OG	2.32	0.42
2:2:67:C:H2'	2:2:68:G:C8	2.54	0.42
2:2:211:G:H3'	2:2:212:G:H4'	2.01	0.42
2:2:982:U:H5''	44:s:6:MET:HE2	1.99	0.42
2:2:1149:C:O2'	2:2:1280:A:N1	2.47	0.42
3:3:6:G:H2'	3:3:7:G:C8	2.54	0.42
13:M:30:SER:HA	13:M:133:LYS:HE2	2.00	0.42
33:h:151:VAL:HG12	33:h:200:VAL:HG23	2.01	0.42
46:u:44:SER:N	46:u:47:GLU:OE2	2.42	0.42
1:1:479:A:N3	1:1:481:G:H5''	2.34	0.42
1:1:1091:G:H1'	31:f:12:ARG:HH22	1.84	0.42
1:1:1590:A:H2'	1:1:1591:A:C8	2.54	0.42
2:2:252:U:O4	2:2:253:A:N6	2.50	0.42
2:2:514:C:H2'	2:2:515:G:C8	2.55	0.42
2:2:1299:A:O2'	2:2:1301:U:O4'	2.27	0.42
3:3:13:G:O2'	3:3:15:A:O5'	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:F:152:ARG:HB2	8:F:162:VAL:H	1.84	0.42
35:j:12:GLN:HE21	35:j:117:VAL:HA	1.83	0.42
42:q:57:LEU:HD12	42:q:61:PHE:HB2	2.00	0.42
54:A:319:ASP:OD1	54:A:320:HIS:N	2.53	0.42
1:1:1297:C:O2'	1:1:1302:A:N1	2.52	0.42
1:1:1682:G:P	1:1:1699:G:H22	2.42	0.42
1:1:2087:G:H2'	1:1:2088:A:C8	2.52	0.42
1:1:2284:A:H1'	1:1:2325:G:C2	2.54	0.42
1:1:2451:A:N1	53:H:78:PHE:CE1	2.82	0.42
1:1:2693:G:H2'	1:1:2694:G:H8	1.83	0.42
2:2:71:A:H61	2:2:99:C:H1'	1.83	0.42
2:2:338:A:H2	2:2:351:G:H22	1.66	0.42
2:2:1032:G:O6	2:2:1034:G:O2'	2.29	0.42
2:2:1308:U:H2'	2:2:1309:G:H8	1.84	0.42
2:2:1437:A:OP1	50:y:29:ARG:NH2	2.39	0.42
8:F:60:ASP:OD1	8:F:60:ASP:N	2.52	0.42
10:J:81:ILE:HD12	10:J:81:ILE:HA	1.92	0.42
18:R:38:VAL:HG13	18:R:54:VAL:HB	2.02	0.42
24:X:12:PRO:HB3	24:X:30:LEU:HD23	2.00	0.42
33:h:29:PHE:HZ	44:s:93:ILE:HG23	1.84	0.42
40:o:6:ILE:HB	40:o:76:ILE:HB	1.99	0.42
43:r:22:ILE:HB	43:r:25:VAL:HG22	2.01	0.42
1:1:155:A:H2'	1:1:156:A:C8	2.55	0.42
1:1:1174:U:H3'	1:1:1175:A:O4'	2.20	0.42
1:1:1464:G:H2'	1:1:1465:G:H8	1.85	0.42
1:1:2503:2MA:O2'	1:1:2505:G:OP2	2.25	0.42
1:1:2520:C:H2'	1:1:2521:C:H6	1.84	0.42
2:2:110:C:O2'	46:u:25:ARG:O	2.37	0.42
2:2:514:C:H2'	2:2:515:G:H8	1.84	0.42
2:2:1499:A:H2'	2:2:1500:A:H8	1.84	0.42
6:D:6:LYS:O	6:D:9:GLN:NE2	2.52	0.42
11:K:66:LYS:N	11:K:82:ASN:OD1	2.44	0.42
24:X:70:GLU:O	24:X:74:ARG:N	2.43	0.42
41:p:112:ASP:HB2	51:z:19:PHE:HE1	1.85	0.42
1:1:549:G:H8	1:1:549:G:O5'	2.02	0.42
1:1:1091:G:N1	1:1:1101:U:O2	2.52	0.42
1:1:2060:A:N7	6:D:69:ARG:NH2	2.59	0.42
3:3:78:A:H62	3:3:98:G:H21	1.67	0.42
13:M:74:THR:HG22	13:M:89:VAL:HA	2.02	0.42
33:h:113:ALA:HB1	33:h:200:VAL:HG13	2.01	0.42
53:H:78:PHE:CD1	53:H:78:PHE:C	2.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:367:G:H3'	1:1:368:A:H8	1.85	0.42
1:1:759:G:H2'	1:1:760:G:C8	2.55	0.42
1:1:1434:A:H2'	1:1:1435:G:C8	2.54	0.42
2:2:213:G:OP2	2:2:214:C:N4	2.39	0.42
2:2:1181:G:H1'	2:2:1182:G:C4	2.55	0.42
2:2:1338:G:N3	52:5:41:C:O2'	2.53	0.42
7:E:52:ASN:HB2	7:E:150:ARG:HH21	1.84	0.42
32:g:73:LYS:HE2	32:g:165:ASP:HB2	2.01	0.42
34:i:170:TRP:CD2	34:i:186:PRO:HB3	2.54	0.42
51:z:4:ILE:HG23	51:z:18:ARG:HE	1.85	0.42
1:1:563:A:OP2	18:R:79:ARG:NH2	2.37	0.42
1:1:602:A:O2'	1:1:604:G:O2'	2.38	0.42
1:1:926:G:H2'	1:1:927:A:C8	2.55	0.42
2:2:1241:G:H2'	2:2:1242:G:H8	1.85	0.42
2:2:1328:C:HO2'	43:r:29:ARG:HH11	1.63	0.42
2:2:1368:A:H5'	39:n:114:LYS:HB3	2.01	0.42
2:2:1386:G:H2'	2:2:1387:G:C8	2.54	0.42
7:E:23:ASN:N	7:E:27:GLN:OE1	2.52	0.42
43:r:80:LEU:HD23	43:r:83:LEU:HD12	2.00	0.42
52:5:63:G:H2'	52:5:64:G:H8	1.84	0.42
1:1:18:U:P	17:Q:30:ARG:HH22	2.43	0.42
1:1:283:G:N2	1:1:358:U:H1'	2.34	0.42
1:1:654:A:O4'	1:1:655:A:OP1	2.37	0.42
1:1:807:U:H2'	1:1:808:G:C8	2.51	0.42
1:1:856:G:H1	1:1:921:C:H42	1.68	0.42
1:1:984:A:H2'	1:1:984:A:N3	2.35	0.42
1:1:1058:U:O4	1:1:1080:A:N6	2.51	0.42
1:1:1436:G:O2'	1:1:1477:A:N3	2.52	0.42
1:1:1466:U:H5''	1:1:1467:U:H5'	2.02	0.42
1:1:1689:A:OP2	1:1:1698:A:N6	2.34	0.42
1:1:1783:A:O2'	1:1:2607:G:O2'	2.29	0.42
1:1:2303:G:O2'	7:E:121:SER:OG	2.34	0.42
1:1:2559:C:H2'	1:1:2560:A:C8	2.54	0.42
2:2:25:C:H2'	2:2:26:A:H8	1.84	0.42
2:2:1162:C:H2'	2:2:1163:A:C8	2.54	0.42
2:2:1525:G:H2'	2:2:1526:G:C8	2.55	0.42
3:3:31:C:H1'	3:3:53:A:N1	2.35	0.42
7:E:65:PRO:HB2	7:E:87:CYS:HB2	2.01	0.42
8:F:156:PRO:O	8:F:172:LYS:N	2.53	0.42
20:T:37:ASP:N	20:T:37:ASP:OD1	2.52	0.42
25:Y:21:LEU:HD12	25:Y:46:VAL:HG13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:g:162:PHE:HA	32:g:184:PHE:HB2	2.01	0.42
34:i:86:THR:O	34:i:90:LEU:N	2.52	0.42
44:s:63:ARG:HB3	44:s:68:GLY:HA2	2.01	0.42
50:y:61:GLN:HB3	50:y:66:LEU:HB3	2.00	0.42
1:1:376:G:H2'	1:1:377:G:C8	2.55	0.42
1:1:510:C:H5'	1:1:510:C:C6	2.55	0.42
1:1:879:G:H1	1:1:899:A:H1'	1.85	0.42
1:1:1035:U:H2'	1:1:1036:G:C8	2.55	0.42
1:1:2685:G:H2'	1:1:2686:G:H8	1.85	0.42
2:2:1:A:O5'	2:2:3:A:N6	2.53	0.42
2:2:1093:A:N3	2:2:1109:C:O2'	2.41	0.42
2:2:1156:G:O2'	2:2:1180:A:N1	2.40	0.42
5:C:46:ARG:HB3	5:C:84:LEU:HD12	2.01	0.42
7:E:175:PHE:HA	7:E:176:PRO:HD3	1.84	0.42
11:K:61:VAL:HB	11:K:87:LEU:HD11	2.02	0.42
32:g:19:GLN:HB2	32:g:22:TYR:CD2	2.55	0.42
33:h:116:VAL:O	33:h:119:SER:OG	2.37	0.42
36:k:29:ILE:HD12	36:k:64:VAL:HG11	2.00	0.42
39:n:21:ILE:HD12	39:n:61:LEU:HD13	2.01	0.42
45:t:26:GLU:OE1	45:t:77:ARG:NH2	2.53	0.42
46:u:45:GLU:OE2	46:u:46:LYS:NZ	2.53	0.42
51:z:18:ARG:O	51:z:22:SER:N	2.47	0.42
54:A:172:TYR:CE2	54:A:207:PRO:HG3	2.55	0.42
1:1:151:C:H2'	1:1:152:A:H8	1.85	0.41
1:1:1166:G:H1	1:1:1183:U:H3	1.68	0.41
1:1:1449:G:H2'	1:1:1450:G:H8	1.85	0.41
1:1:1934:C:H2'	1:1:1935:G:C8	2.55	0.41
1:1:2544:G:H8	1:1:2544:G:O5'	2.02	0.41
2:2:343:U:O2'	2:2:346:G:O6	2.30	0.41
2:2:363:A:H5'	42:q:31:ARG:HB2	2.01	0.41
2:2:407:U:H5''	34:i:112:ALA:HB1	2.01	0.41
2:2:526:C:C4	2:2:527:7MG:H1'	2.55	0.41
2:2:986:U:H4'	49:x:55:ARG:HG3	2.01	0.41
24:X:43:GLU:OE1	24:X:45:ARG:NE	2.38	0.41
34:i:122:ALA:C	34:i:146:ARG:CG	2.94	0.41
40:o:42:LEU:HD12	40:o:71:LEU:HB3	2.01	0.41
52:5:17:U:H4'	52:5:18:G:H5''	2.02	0.41
52:5:66:C:H2'	52:5:67:C:O4'	2.20	0.41
1:1:458:G:O2'	1:1:469:G:O6	2.29	0.41
1:1:2313:C:H2'	1:1:2314:A:C8	2.55	0.41
2:2:72:A:C5	2:2:73:C:H1'	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:75:G:H2'	2:2:76:G:H8	1.84	0.41
2:2:214:C:H2'	2:2:215:C:C6	2.56	0.41
2:2:322:C:OP2	2:2:328:C:N4	2.51	0.41
2:2:1350:A:OP1	39:n:123:ARG:NH1	2.53	0.41
3:3:92:C:H2'	3:3:93:C:C6	2.55	0.41
33:h:141:ALA:HA	33:h:144:LEU:HB2	2.02	0.41
52:5:7:U:O2'	52:5:21:A:N1	2.44	0.41
54:A:106:ASP:OD1	54:A:109:ARG:NE	2.42	0.41
1:1:161:A:HO2'	1:1:2207:C:HO2'	1.69	0.41
1:1:742:A:H2'	1:1:743:A:C8	2.55	0.41
1:1:767:U:H2'	1:1:768:G:H8	1.86	0.41
1:1:1464:G:H2'	1:1:1465:G:C8	2.55	0.41
1:1:1509:A:H1'	1:1:1510:G:N9	2.35	0.41
1:1:1586:A:H3'	1:1:1587:G:C8	2.55	0.41
1:1:1980:G:O2'	1:1:1982:U:OP2	2.36	0.41
1:1:2293:G:H5''	15:O:94:ARG:HH12	1.86	0.41
1:1:2315:G:H2'	1:1:2316:G:H8	1.85	0.41
1:1:2801:G:H2'	1:1:2802:G:H8	1.86	0.41
2:2:123:U:O2'	2:2:290:C:O2'	2.32	0.41
2:2:219:U:H2'	2:2:220:G:C8	2.55	0.41
2:2:591:U:H2'	2:2:592:G:C8	2.56	0.41
2:2:744:C:H2'	2:2:745:G:C8	2.54	0.41
34:i:95:GLU:HA	34:i:100:ASN:ND2	2.35	0.41
1:1:597:G:O2'	12:L:11:GLY:O	2.32	0.41
1:1:2395:C:OP1	12:L:63:LYS:NZ	2.42	0.41
1:1:2861:U:H2'	1:1:2862:G:H8	1.84	0.41
2:2:120:A:O2'	2:2:122:G:N7	2.46	0.41
2:2:335:C:H2'	2:2:336:A:C8	2.55	0.41
2:2:1440:U:H2'	2:2:1441:A:C8	2.56	0.41
3:3:116:G:H2'	3:3:117:G:C8	2.56	0.41
10:J:60:ASP:OD1	10:J:60:ASP:N	2.52	0.41
18:R:77:PHE:HD1	18:R:84:ARG:HB3	1.85	0.41
41:p:88:GLY:H	41:p:114:THR:HG22	1.85	0.41
42:q:36:ARG:HA	42:q:76:GLU:HG3	2.01	0.41
44:s:24:ARG:HD2	44:s:55:SER:HB3	2.02	0.41
54:A:210:PRO:HD2	54:A:213:GLU:HB2	2.03	0.41
1:1:58:G:O2'	1:1:73:A:N1	2.36	0.41
1:1:441:U:H2'	1:1:442:G:C8	2.56	0.41
1:1:1483:G:O4'	1:1:1510:G:N2	2.45	0.41
1:1:1533:C:H2'	1:1:1535:A:H2	1.85	0.41
1:1:2452:C:H2'	1:1:2453:A:C8	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2698:U:H2'	1:1:2699:C:C6	2.55	0.41
2:2:376:G:H2'	2:2:377:G:C8	2.55	0.41
9:G:51:ARG:HA	9:G:54:LEU:HB3	2.03	0.41
34:i:173:VAL:HA	34:i:180:GLY:HA2	2.03	0.41
45:t:33:THR:HG21	45:t:85:LEU:HD12	2.03	0.41
1:1:20:C:H2'	1:1:21:A:H8	1.86	0.41
1:1:287:G:H2'	1:1:288:U:C6	2.56	0.41
1:1:571:U:H2'	18:R:80:ARG:HH12	1.85	0.41
1:1:647:G:C8	1:1:647:G:C5'	2.88	0.41
1:1:762:U:N3	1:1:1431:A:OP1	2.51	0.41
1:1:893:C:H2'	1:1:894:U:C5	2.55	0.41
1:1:1298:C:H42	1:1:1642:G:H1	1.67	0.41
1:1:1329:U:O2'	1:1:1330:C:H5'	2.21	0.41
1:1:1438:U:H2'	1:1:1439:A:C8	2.56	0.41
1:1:1682:G:C2	1:1:1757:A:H1'	2.55	0.41
1:1:1722:A:H62	1:1:1738:G:H1'	1.86	0.41
1:1:2246:G:H2'	1:1:2247:A:C8	2.56	0.41
1:1:2286:G:H4'	1:1:2287:A:O4'	2.21	0.41
1:1:2809:A:OP2	1:1:2890:G:N1	2.36	0.41
2:2:962:C:H2'	2:2:963:G:C8	2.56	0.41
2:2:1025:U:H4'	2:2:1026:G:C8	2.55	0.41
2:2:1321:U:O3'	49:x:78:ARG:NH2	2.54	0.41
10:J:32:LEU:HD22	10:J:54:ILE:HG21	2.02	0.41
32:g:32:PHE:N	32:g:40:ILE:O	2.53	0.41
35:j:83:HIS:NE2	35:j:147:MET:HG3	2.36	0.41
38:m:79:SER:HB2	38:m:85:ILE:HB	2.02	0.41
39:n:12:ARG:HH22	39:n:109:ARG:NH2	2.19	0.41
52:5:5:G:H2'	52:5:6:G:C8	2.55	0.41
52:5:55:PSU:HN3	52:5:57:A:H3'	1.85	0.41
54:A:244:ILE:HD11	54:A:263:GLN:HG3	2.03	0.41
1:1:1350:C:C5	1:1:1351:C:N4	2.86	0.41
1:1:1492:G:H1	1:1:1498:C:H42	1.67	0.41
1:1:1965:C:H5''	1:1:1966:A:H2'	2.03	0.41
1:1:1969:A:O2'	1:1:1972:G:N3	2.51	0.41
1:1:2615:U:H2'	1:1:2616:C:C6	2.56	0.41
1:1:2735:G:H2'	1:1:2736:A:H8	1.86	0.41
2:2:409:U:OP1	34:i:23:SER:HB3	2.21	0.41
2:2:1308:U:H2'	2:2:1309:G:C8	2.56	0.41
7:E:34:ILE:HG12	7:E:156:ILE:HG12	2.02	0.41
27:b:40:ARG:HG3	27:b:41:HIS:ND1	2.35	0.41
28:c:15:ALA:HB2	28:c:47:VAL:HG21	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:n:6:TYR:O	39:n:21:ILE:N	2.53	0.41
54:A:127:PHE:HA	54:A:130:ASP:HB2	2.02	0.41
54:A:311:ASN:HB3	54:A:316:ARG:HB3	2.01	0.41
1:1:300:A:H2'	1:1:334:C:H1'	2.02	0.41
1:1:581:C:H2'	1:1:582:A:C8	2.56	0.41
1:1:680:C:H2'	1:1:681:G:C8	2.55	0.41
1:1:1341:G:OP2	1:1:1394:U:O2'	2.31	0.41
1:1:1350:C:C2	1:1:1351:C:C4	3.09	0.41
1:1:1771:C:H2'	1:1:1772:A:C8	2.55	0.41
1:1:1864:U:H3	1:1:1878:G:H1	1.68	0.41
1:1:2315:G:H2'	1:1:2316:G:C8	2.56	0.41
2:2:555:U:H2'	2:2:556:C:C6	2.56	0.41
2:2:662:U:H2'	2:2:663:A:C8	2.55	0.41
2:2:1056:U:H2'	2:2:1057:G:H8	1.85	0.41
2:2:1260:G:H21	2:2:1275:A:H8	1.69	0.41
33:h:142:MET:HE2	33:h:170:GLU:HG2	2.03	0.41
36:k:81:ASN:HB3	36:k:84:VAL:HG12	2.03	0.41
38:m:34:VAL:O	38:m:38:ASN:N	2.53	0.41
39:n:50:GLN:OE1	39:n:80:ARG:NH1	2.52	0.41
39:n:63:LEU:HD13	39:n:65:ILE:HD12	2.03	0.41
41:p:46:THR:OG1	41:p:49:GLY:N	2.45	0.41
1:1:544:C:H3'	1:1:545:U:O4'	2.21	0.41
1:1:548:G:H3'	1:1:549:G:C8	2.56	0.41
1:1:631:A:N3	1:1:2415:G:O2'	2.51	0.41
1:1:748:G:OP1	19:S:88:ARG:NH1	2.38	0.41
1:1:871:U:H2'	1:1:872:U:C6	2.56	0.41
1:1:1041:G:H2'	1:1:1042:G:C8	2.56	0.41
1:1:1599:U:OP1	20:T:39:THR:OG1	2.34	0.41
1:1:2304:G:O2'	7:E:130:MET:O	2.27	0.41
1:1:2340:A:H5'	3:3:41:G:H21	1.86	0.41
1:1:2467:C:O2	13:M:123:LYS:NZ	2.53	0.41
2:2:25:C:H2'	2:2:26:A:C8	2.55	0.41
2:2:62:U:O2	2:2:105:G:N2	2.37	0.41
2:2:707:U:H4'	41:p:22:HIS:CD2	2.56	0.41
2:2:1094:G:O2'	2:2:1108:G:N1	2.54	0.41
2:2:1193:G:N7	33:h:3:GLN:NE2	2.55	0.41
2:2:1297:G:N2	2:2:1298:U:O4	2.32	0.41
2:2:1461:G:H2'	2:2:1462:C:C6	2.56	0.41
3:3:85:G:H2'	3:3:86:G:H8	1.85	0.41
3:3:114:C:H2'	3:3:115:A:H8	1.86	0.41
5:C:45:TYR:OH	5:C:81:GLU:OE2	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:187:LEU:HD12	5:C:187:LEU:HA	1.93	0.41
7:E:66:LEU:N	7:E:88:LYS:O	2.50	0.41
9:G:55:GLU:HA	9:G:58:LEU:HB2	2.02	0.41
9:G:81:ALA:HB1	9:G:149:GLU:HB2	2.03	0.41
10:J:49:ASP:CG	10:J:121:LYS:HZ3	2.29	0.41
14:N:35:LYS:HE3	14:N:100:CYS:SG	2.61	0.41
21:U:13:VAL:HG21	21:U:39:ILE:HG21	2.03	0.41
21:U:43:LYS:HA	21:U:60:GLU:HA	2.03	0.41
25:Y:21:LEU:O	25:Y:25:GLN:N	2.46	0.41
27:b:38:HIS:ND1	27:b:39:LEU:O	2.51	0.41
35:j:152:MET:O	35:j:156:LYS:N	2.53	0.41
36:k:75:GLU:HG2	36:k:89:VAL:HG21	2.02	0.41
40:o:29:ALA:HB1	40:o:83:THR:HG23	2.03	0.41
44:s:64:CYS:HB3	44:s:69:ARG:H	1.86	0.41
54:A:253:ILE:HD12	54:A:278:ILE:HD13	2.02	0.41
1:1:1141:U:H4'	1:1:1142:A:O4'	2.21	0.41
1:1:1565:C:O2'	1:1:1567:G:N7	2.34	0.41
1:1:1867:G:H2'	1:1:1868:C:C6	2.56	0.41
1:1:2171:A:H3'	1:1:2174:C:H5	1.84	0.41
1:1:2491:U:H5'	1:1:2570:G:H5''	2.03	0.41
1:1:2848:G:O2'	1:1:2867:G:N2	2.38	0.41
1:1:2880:C:O2'	14:N:90:ARG:NH2	2.51	0.41
2:2:157:U:H2'	2:2:158:G:C8	2.55	0.41
2:2:244:U:O4	2:2:892:A:N6	2.54	0.41
2:2:1152:A:O5'	40:o:72:ARG:NH2	2.54	0.41
2:2:1180:A:OP1	39:n:105:THR:OG1	2.33	0.41
2:2:1256:A:O2'	2:2:1278:G:O6	2.37	0.41
2:2:1275:A:H3'	2:2:1276:G:C8	2.51	0.41
2:2:1303:C:H2'	2:2:1304:G:O4'	2.21	0.41
7:E:4:LEU:HA	7:E:4:LEU:HD23	1.92	0.41
21:U:51:ALA:O	21:U:52:LEU:CB	2.69	0.41
22:V:75:GLN:HB2	22:V:92:VAL:HG23	2.03	0.41
29:d:25:LYS:HE3	29:d:25:LYS:HB3	1.95	0.41
39:n:97:GLU:O	39:n:101:ALA:N	2.53	0.41
41:p:96:THR:O	41:p:100:LEU:N	2.54	0.41
42:q:89:0TD:H8	42:q:89:0TD:H4	1.82	0.41
44:s:49:GLN:OE1	49:x:13:LEU:N	2.53	0.41
46:u:44:SER:OG	46:u:45:GLU:N	2.55	0.41
52:5:13:A:H3'	52:5:14:G:H8	1.86	0.41
1:1:2029:G:N1	1:1:2033:A:OP2	2.38	0.40
1:1:2123:G:N2	1:1:2176:A:O2'	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:2619:C:H5''	5:C:157:LYS:HD2	2.04	0.40
1:1:2870:C:H2'	1:1:2871:U:O4'	2.21	0.40
2:2:83:C:H1'	2:2:86:G:H22	1.86	0.40
2:2:986:U:O2'	49:x:55:ARG:O	2.38	0.40
2:2:1376:U:H2'	2:2:1377:A:C8	2.54	0.40
2:2:1530:G:H2'	2:2:1531:A:H8	1.86	0.40
13:M:108:VAL:HB	13:M:112:LEU:HD23	2.02	0.40
33:h:20:SER:OG	33:h:40:ARG:NH2	2.38	0.40
34:i:69:GLU:OE2	34:i:204:TYR:OH	2.28	0.40
47:v:68:SER:OG	47:v:69:LYS:N	2.54	0.40
54:A:138:TYR:OH	54:A:338:ASP:OD1	2.32	0.40
54:A:307:ASN:O	54:A:320:HIS:N	2.53	0.40
1:1:605:G:OP1	6:D:99:LYS:NZ	2.54	0.40
1:1:621:A:OP2	12:L:99:ASN:ND2	2.54	0.40
1:1:714:U:H1'	1:1:717:C:H5	1.86	0.40
1:1:935:C:H2'	1:1:936:A:H8	1.86	0.40
1:1:1212:G:O2'	1:1:1236:G:N1	2.40	0.40
1:1:1351:C:H6	1:1:1351:C:O5'	2.04	0.40
1:1:2029:G:O5'	1:1:2029:G:H8	2.04	0.40
1:1:2036:C:H2'	1:1:2037:A:H8	1.85	0.40
1:1:2810:A:H62	1:1:2890:G:H21	1.70	0.40
2:2:376:G:H5''	46:u:5:ARG:HD2	2.03	0.40
2:2:1125:U:H5'	40:o:7:ARG:HH12	1.86	0.40
2:2:1507:A:H2'	2:2:1508:A:C8	2.56	0.40
2:2:1513:A:H2'	2:2:1514:G:C8	2.56	0.40
9:G:116:ARG:HG3	9:G:133:GLN:HB2	2.03	0.40
10:J:49:ASP:HB2	10:J:114:LEU:HD11	2.03	0.40
12:L:29:LYS:HG3	12:L:30:THR:HG23	2.02	0.40
13:M:69:PRO:HA	13:M:94:ALA:HB2	2.02	0.40
22:V:6:ALA:N	22:V:64:VAL:O	2.54	0.40
33:h:60:PRO:O	33:h:63:SER:OG	2.37	0.40
44:s:99:ALA:HB1	44:s:101:TRP:HZ3	1.85	0.40
46:u:40:ASN:HB3	46:u:43:ALA:HB2	2.04	0.40
54:A:183:ARG:O	54:A:310:TYR:N	2.46	0.40
54:A:343:PRO:HA	54:A:346:GLN:HB2	2.01	0.40
1:1:379:G:H1	1:1:395:U:H3	1.69	0.40
1:1:393:C:H2'	1:1:394:C:C6	2.55	0.40
1:1:523:C:H2'	1:1:524:G:H8	1.86	0.40
1:1:668:A:H2'	1:1:670:A:H62	1.86	0.40
1:1:953:G:H2'	1:1:954:G:H8	1.86	0.40
1:1:1318:U:H2'	1:1:1319:C:C6	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1340:U:OP1	20:T:19:LYS:NZ	2.48	0.40
1:1:2393:U:H2'	1:1:2394:C:H6	1.86	0.40
1:1:2615:U:C2	27:b:4:GLN:HA	2.57	0.40
2:2:55:A:H61	2:2:357:G:H1'	1.85	0.40
2:2:377:G:H2'	2:2:378:G:H8	1.85	0.40
8:F:105:LEU:HB2	8:F:113:VAL:HB	2.02	0.40
11:K:38:ILE:HD11	11:K:112:PHE:HZ	1.87	0.40
33:h:22:TRP:HA	40:o:94:ALA:HB1	2.04	0.40
1:1:409:G:H2'	1:1:410:G:C8	2.56	0.40
1:1:639:U:H2'	1:1:640:C:C6	2.56	0.40
1:1:849:A:H61	1:1:928:A:H61	1.70	0.40
1:1:897:C:H2'	1:1:898:C:O4'	2.22	0.40
1:1:946:C:H2'	1:1:947:A:H8	1.85	0.40
1:1:1095:A:O2'	1:1:1096:A:O5'	2.36	0.40
1:1:1381:G:H2'	1:1:1382:G:C8	2.56	0.40
1:1:2210:U:H4'	1:1:2211:A:H2'	2.03	0.40
2:2:1002:G:H2'	2:2:1003:G:H8	1.86	0.40
2:2:1178:G:HO2'	2:2:1180:A:H62	1.64	0.40
7:E:158:THR:HG1	7:E:159:THR:H	1.67	0.40
33:h:65:ARG:NE	33:h:100:GLN:OE1	2.49	0.40
34:i:9:LEU:HD11	34:i:31:LYS:HE3	2.03	0.40
34:i:139:PRO:HA	34:i:182:PHE:HD2	1.86	0.40
38:m:89:LYS:NZ	38:m:90:ASP:OD2	2.53	0.40
41:p:94:GLU:OE2	41:p:98:ARG:NH2	2.49	0.40
44:s:37:SER:HB2	44:s:40:ASP:HB2	2.02	0.40
54:A:253:ILE:HD13	54:A:277:ARG:HB3	2.03	0.40
1:1:301:G:H5'	1:1:334:C:H2'	2.04	0.40
1:1:613:A:N3	1:1:614:A:C8	2.90	0.40
1:1:875:G:N2	1:1:903:C:H1'	2.37	0.40
1:1:1158:C:H5''	26:Z:31:ARG:HG3	2.03	0.40
1:1:1165:A:H2'	1:1:1166:G:H8	1.86	0.40
1:1:1386:C:H2'	1:1:1387:A:H8	1.85	0.40
1:1:2461:A:N6	1:1:2488:G:O6	2.55	0.40
2:2:424:G:H2'	2:2:425:G:C8	2.56	0.40
11:K:121:GLU:OE1	16:P:65:SER:OG	2.36	0.40
19:S:81:SER:HB3	19:S:97:LEU:HD12	2.02	0.40
25:Y:18:LEU:HB2	25:Y:53:VAL:HG11	2.02	0.40
35:j:72:ILE:HD12	35:j:72:ILE:HA	1.98	0.40
35:j:103:THR:N	35:j:122:ASN:OD1	2.54	0.40
52:5:63:G:H2'	52:5:64:G:C8	2.56	0.40
54:A:223:ARG:HB2	54:A:249:LEU:HD21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	
4	B	269/271 (99%)	247 (92%)	22 (8%)	0	100	100
5	C	207/209 (99%)	194 (94%)	12 (6%)	1 (0%)	24	57
6	D	199/201 (99%)	191 (96%)	8 (4%)	0	100	100
7	E	175/177 (99%)	155 (89%)	19 (11%)	1 (1%)	21	54
8	F	173/175 (99%)	161 (93%)	12 (7%)	0	100	100
9	G	147/149 (99%)	133 (90%)	14 (10%)	0	100	100
10	J	140/142 (99%)	136 (97%)	4 (3%)	0	100	100
11	K	121/123 (98%)	111 (92%)	10 (8%)	0	100	100
12	L	142/144 (99%)	135 (95%)	7 (5%)	0	100	100
13	M	134/136 (98%)	127 (95%)	7 (5%)	0	100	100
14	N	117/119 (98%)	109 (93%)	8 (7%)	0	100	100
15	O	114/116 (98%)	111 (97%)	3 (3%)	0	100	100
16	P	112/114 (98%)	108 (96%)	4 (4%)	0	100	100
17	Q	115/117 (98%)	113 (98%)	2 (2%)	0	100	100
18	R	101/103 (98%)	92 (91%)	7 (7%)	2 (2%)	6	32
19	S	108/110 (98%)	101 (94%)	7 (6%)	0	100	100
20	T	92/94 (98%)	89 (97%)	3 (3%)	0	100	100
21	U	101/103 (98%)	89 (88%)	12 (12%)	0	100	100
22	V	92/94 (98%)	87 (95%)	5 (5%)	0	100	100
23	W	74/76 (97%)	68 (92%)	6 (8%)	0	100	100
24	X	75/77 (97%)	71 (95%)	4 (5%)	0	100	100
25	Y	60/62 (97%)	58 (97%)	2 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	Z	56/58 (97%)	53 (95%)	3 (5%)	0	100	100
27	b	54/56 (96%)	51 (94%)	3 (6%)	0	100	100
28	c	50/52 (96%)	49 (98%)	1 (2%)	0	100	100
29	d	44/46 (96%)	42 (96%)	2 (4%)	0	100	100
30	e	62/64 (97%)	57 (92%)	4 (6%)	1 (2%)	7	36
31	f	36/38 (95%)	36 (100%)	0	0	100	100
32	g	223/225 (99%)	201 (90%)	22 (10%)	0	100	100
33	h	206/208 (99%)	192 (93%)	14 (7%)	0	100	100
34	i	203/205 (99%)	188 (93%)	13 (6%)	2 (1%)	12	44
35	j	154/156 (99%)	140 (91%)	14 (9%)	0	100	100
36	k	102/104 (98%)	100 (98%)	2 (2%)	0	100	100
37	l	149/151 (99%)	144 (97%)	5 (3%)	0	100	100
38	m	127/129 (98%)	119 (94%)	8 (6%)	0	100	100
39	n	125/127 (98%)	114 (91%)	11 (9%)	0	100	100
40	o	97/99 (98%)	90 (93%)	7 (7%)	0	100	100
41	p	115/117 (98%)	110 (96%)	5 (4%)	0	100	100
42	q	120/123 (98%)	109 (91%)	11 (9%)	0	100	100
43	r	114/116 (98%)	108 (95%)	6 (5%)	0	100	100
44	s	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
45	t	86/88 (98%)	80 (93%)	6 (7%)	0	100	100
46	u	80/82 (98%)	74 (92%)	6 (8%)	0	100	100
47	v	78/80 (98%)	69 (88%)	9 (12%)	0	100	100
48	w	64/66 (97%)	63 (98%)	1 (2%)	0	100	100
49	x	81/83 (98%)	79 (98%)	2 (2%)	0	100	100
50	y	84/86 (98%)	81 (96%)	3 (4%)	0	100	100
51	z	68/70 (97%)	66 (97%)	2 (3%)	0	100	100
53	H	1/3 (33%)	1 (100%)	0	0	100	100
54	A	242/257 (94%)	236 (98%)	6 (2%)	0	100	100
All	All	5787/5901 (98%)	5433 (94%)	347 (6%)	7 (0%)	49	79

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
18	R	54	VAL
34	i	146	ARG
18	R	52	PRO
34	i	148	LYS
7	E	123	ASP
30	e	32	ILE
5	C	152	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 PDB entries followed by that with respect to all EM entries.

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	
4	B	216/216 (100%)	216 (100%)	0	100	100
5	C	164/164 (100%)	163 (99%)	1 (1%)	78	79
6	D	165/165 (100%)	163 (99%)	2 (1%)	63	73
7	E	148/148 (100%)	148 (100%)	0	100	100
8	F	136/136 (100%)	135 (99%)	1 (1%)	76	78
9	G	114/114 (100%)	114 (100%)	0	100	100
10	J	116/116 (100%)	116 (100%)	0	100	100
11	K	104/104 (100%)	104 (100%)	0	100	100
12	L	103/103 (100%)	103 (100%)	0	100	100
13	M	109/109 (100%)	109 (100%)	0	100	100
14	N	99/99 (100%)	99 (100%)	0	100	100
15	O	86/86 (100%)	86 (100%)	0	100	100
16	P	99/99 (100%)	98 (99%)	1 (1%)	68	75
17	Q	89/89 (100%)	89 (100%)	0	100	100
18	R	84/84 (100%)	81 (96%)	3 (4%)	31	57
19	S	93/93 (100%)	93 (100%)	0	100	100
20	T	81/81 (100%)	81 (100%)	0	100	100
21	U	84/84 (100%)	83 (99%)	1 (1%)	63	73
22	V	78/78 (100%)	77 (99%)	1 (1%)	61	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
23	W	58/58 (100%)	58 (100%)	0	100	100
24	X	67/67 (100%)	67 (100%)	0	100	100
25	Y	54/54 (100%)	54 (100%)	0	100	100
26	Z	48/48 (100%)	48 (100%)	0	100	100
27	b	47/47 (100%)	47 (100%)	0	100	100
28	c	47/47 (100%)	47 (100%)	0	100	100
29	d	38/38 (100%)	38 (100%)	0	100	100
30	e	51/51 (100%)	51 (100%)	0	100	100
31	f	34/34 (100%)	33 (97%)	1 (3%)	37	60
32	g	187/187 (100%)	187 (100%)	0	100	100
33	h	171/171 (100%)	170 (99%)	1 (1%)	78	79
34	i	172/172 (100%)	170 (99%)	2 (1%)	63	73
35	j	119/119 (100%)	119 (100%)	0	100	100
36	k	91/91 (100%)	91 (100%)	0	100	100
37	l	124/124 (100%)	124 (100%)	0	100	100
38	m	104/104 (100%)	104 (100%)	0	100	100
39	n	105/105 (100%)	105 (100%)	0	100	100
40	o	86/86 (100%)	86 (100%)	0	100	100
41	p	90/90 (100%)	90 (100%)	0	100	100
42	q	102/102 (100%)	102 (100%)	0	100	100
43	r	94/94 (100%)	94 (100%)	0	100	100
44	s	83/83 (100%)	83 (100%)	0	100	100
45	t	76/76 (100%)	76 (100%)	0	100	100
46	u	65/65 (100%)	65 (100%)	0	100	100
47	v	74/74 (100%)	74 (100%)	0	100	100
48	w	57/57 (100%)	57 (100%)	0	100	100
49	x	72/72 (100%)	72 (100%)	0	100	100
50	y	65/65 (100%)	65 (100%)	0	100	100
51	z	60/60 (100%)	60 (100%)	0	100	100
53	H	2/2 (100%)	0	2 (100%)	0	0
54	A	201/208 (97%)	201 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	4812/4819 (100%)	4796 (100%)	16 (0%)	84 83

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	C	151	THR
6	D	163	ASN
6	D	170	ARG
8	F	11	VAL
16	P	114	LEU
18	R	51	VAL
18	R	54	VAL
18	R	80	ARG
21	U	52	LEU
22	V	77	VAL
31	f	26	ILE
33	h	200	VAL
34	i	146	ARG
34	i	147	GLU
53	H	78	PHE
53	H	79	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
4	B	86	ASN
4	B	134	ASN
4	B	197	ASN
4	B	260	ASN
5	C	49	GLN
5	C	136	ASN
6	D	115	GLN
6	D	136	GLN
6	D	163	ASN
8	F	104	ASN
15	O	29	HIS
15	O	38	GLN
16	P	10	GLN
16	P	12	GLN
16	P	15	GLN
16	P	56	HIS

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Mol	Chain	Res	Type
16	P	66	ASN
17	Q	44	GLN
19	S	7	HIS
19	S	61	ASN
21	U	46	GLN
21	U	54	GLN
22	V	24	ASN
22	V	75	GLN
25	Y	15	ASN
25	Y	45	GLN
28	c	26	ASN
30	e	31	HIS
31	f	13	ASN
32	g	18	HIS
32	g	39	HIS
33	h	41	GLN
33	h	190	HIS
34	i	89	ASN
34	i	198	HIS
35	j	97	GLN
35	j	132	ASN
36	k	46	GLN
36	k	55	HIS
36	k	58	HIS
36	k	63	ASN
37	l	28	ASN
37	l	68	ASN
37	l	97	ASN
37	l	130	ASN
39	n	31	ASN
41	p	22	HIS
41	p	28	ASN
41	p	40	ASN
42	q	5	ASN
42	q	6	GLN
44	s	4	GLN
45	t	20	ASN
45	t	37	ASN
45	t	40	GLN
46	u	9	HIS
46	u	79	ASN
49	x	83	HIS

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Mol	Chain	Res	Type
50	y	20	HIS
50	y	52	ASN
50	y	61	GLN
54	A	307	ASN
54	A	352	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	1	2902/2903 (99%)	693 (23%)	15 (0%)
2	2	1532/1534 (99%)	363 (23%)	6 (0%)
3	3	119/120 (99%)	30 (25%)	0
52	5	75/76 (98%)	24 (32%)	1 (1%)
55	4	8/9 (88%)	2 (25%)	0
All	All	4636/4642 (99%)	1112 (23%)	22 (0%)

All (1112) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	1	2	G
1	1	10	A
1	1	15	G
1	1	34	U
1	1	35	G
1	1	36	G
1	1	46	G
1	1	54	G
1	1	55	G
1	1	62	U
1	1	63	A
1	1	71	A
1	1	73	A
1	1	74	A
1	1	75	G
1	1	80	G
1	1	85	G
1	1	87	U
1	1	88	G
1	1	91	A
1	1	96	C
1	1	101	A

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Mol	Chain	Res	Type
1	1	102	U
1	1	103	A
1	1	110	G
1	1	118	A
1	1	119	A
1	1	120	U
1	1	125	A
1	1	126	A
1	1	132	G
1	1	136	G
1	1	137	U
1	1	138	U
1	1	139	U
1	1	140	C
1	1	142	A
1	1	160	A
1	1	163	C
1	1	179	C
1	1	181	A
1	1	184	C
1	1	196	A
1	1	199	A
1	1	213	A
1	1	215	G
1	1	216	A
1	1	222	A
1	1	225	C
1	1	228	C
1	1	229	C
1	1	230	G
1	1	239	C
1	1	241	A
1	1	248	G
1	1	250	G
1	1	265	A
1	1	266	G
1	1	268	C
1	1	270	A
1	1	271	G
1	1	272	A
1	1	273	G
1	1	275	C

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Mol	Chain	Res	Type
1	1	276	U
1	1	277	G
1	1	278	A
1	1	279	A
1	1	284	U
1	1	285	G
1	1	286	U
1	1	292	U
1	1	295	G
1	1	311	A
1	1	312	G
1	1	315	G
1	1	323	C
1	1	329	G
1	1	330	A
1	1	345	A
1	1	352	A
1	1	353	C
1	1	354	A
1	1	360	U
1	1	361	G
1	1	367	G
1	1	369	U
1	1	371	A
1	1	372	G
1	1	383	C
1	1	386	G
1	1	387	U
1	1	388	G
1	1	389	G
1	1	391	A
1	1	393	C
1	1	404	A
1	1	405	U
1	1	411	G
1	1	412	A
1	1	424	G
1	1	435	C
1	1	438	G
1	1	448	U
1	1	449	A
1	1	456	C

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Mol	Chain	Res	Type
1	1	457	A
1	1	467	G
1	1	477	A
1	1	479	A
1	1	480	A
1	1	481	G
1	1	489	G
1	1	491	G
1	1	496	G
1	1	501	A
1	1	505	A
1	1	509	C
1	1	510	C
1	1	527	C
1	1	529	A
1	1	530	G
1	1	531	C
1	1	532	A
1	1	533	G
1	1	543	G
1	1	544	C
1	1	545	U
1	1	546	U
1	1	547	A
1	1	548	G
1	1	549	G
1	1	551	G
1	1	563	A
1	1	573	U
1	1	575	A
1	1	586	A
1	1	587	C
1	1	602	A
1	1	603	A
1	1	609	A
1	1	613	A
1	1	614	A
1	1	615	U
1	1	616	A
1	1	621	A
1	1	627	A
1	1	637	A

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Mol	Chain	Res	Type
1	1	645	C
1	1	646	U
1	1	647	G
1	1	648	G
1	1	651	G
1	1	653	U
1	1	654	A
1	1	655	A
1	1	656	G
1	1	659	G
1	1	664	G
1	1	685	A
1	1	686	U
1	1	709	U
1	1	710	U
1	1	717	C
1	1	726	G
1	1	729	G
1	1	730	A
1	1	734	A
1	1	738	G
1	1	747	5MU
1	1	764	A
1	1	765	C
1	1	770	G
1	1	775	G
1	1	776	G
1	1	782	A
1	1	783	A
1	1	784	G
1	1	785	G
1	1	792	A
1	1	797	G
1	1	798	G
1	1	800	A
1	1	805	G
1	1	811	U
1	1	812	C
1	1	819	A
1	1	827	U
1	1	828	U
1	1	831	G

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Mol	Chain	Res	Type
1	1	846	U
1	1	856	G
1	1	858	G
1	1	859	G
1	1	866	A
1	1	869	G
1	1	877	A
1	1	878	A
1	1	879	G
1	1	880	G
1	1	882	G
1	1	884	U
1	1	885	C
1	1	887	A
1	1	889	C
1	1	890	C
1	1	891	G
1	1	893	C
1	1	894	U
1	1	895	U
1	1	896	A
1	1	897	C
1	1	899	A
1	1	907	G
1	1	910	A
1	1	914	G
1	1	915	C
1	1	919	U
1	1	931	U
1	1	932	U
1	1	941	A
1	1	946	C
1	1	957	C
1	1	959	A
1	1	961	C
1	1	962	G
1	1	973	A
1	1	974	G
1	1	983	A
1	1	989	G
1	1	990	A
1	1	995	C

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Mol	Chain	Res	Type
1	1	996	A
1	1	1012	U
1	1	1013	C
1	1	1022	G
1	1	1025	G
1	1	1030	C
1	1	1033	U
1	1	1034	G
1	1	1040	A
1	1	1042	G
1	1	1045	C
1	1	1046	A
1	1	1047	G
1	1	1051	G
1	1	1058	U
1	1	1059	G
1	1	1060	U
1	1	1061	U
1	1	1064	C
1	1	1065	U
1	1	1066	U
1	1	1067	A
1	1	1068	G
1	1	1069	A
1	1	1070	A
1	1	1071	G
1	1	1074	G
1	1	1075	C
1	1	1078	U
1	1	1079	C
1	1	1083	U
1	1	1084	A
1	1	1087	G
1	1	1088	A
1	1	1090	A
1	1	1095	A
1	1	1096	A
1	1	1097	U
1	1	1110	G
1	1	1111	A
1	1	1112	G
1	1	1119	U

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Mol	Chain	Res	Type
1	1	1121	C
1	1	1129	A
1	1	1130	U
1	1	1131	G
1	1	1132	U
1	1	1133	A
1	1	1134	A
1	1	1135	C
1	1	1136	G
1	1	1139	G
1	1	1142	A
1	1	1151	A
1	1	1155	A
1	1	1168	G
1	1	1169	A
1	1	1170	C
1	1	1173	U
1	1	1174	U
1	1	1175	A
1	1	1176	U
1	1	1177	G
1	1	1178	C
1	1	1179	G
1	1	1180	U
1	1	1182	G
1	1	1186	G
1	1	1204	A
1	1	1205	A
1	1	1212	G
1	1	1218	G
1	1	1238	G
1	1	1247	A
1	1	1248	G
1	1	1252	G
1	1	1253	A
1	1	1256	G
1	1	1265	A
1	1	1266	G
1	1	1271	G
1	1	1272	A
1	1	1273	U
1	1	1300	G

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Mol	Chain	Res	Type
1	1	1301	A
1	1	1313	U
1	1	1315	C
1	1	1321	A
1	1	1325	U
1	1	1350	C
1	1	1352	U
1	1	1360	G
1	1	1365	A
1	1	1368	G
1	1	1374	G
1	1	1379	U
1	1	1380	G
1	1	1383	A
1	1	1395	A
1	1	1398	C
1	1	1407	G
1	1	1408	G
1	1	1409	U
1	1	1410	G
1	1	1411	U
1	1	1415	U
1	1	1417	C
1	1	1419	A
1	1	1420	A
1	1	1427	A
1	1	1428	C
1	1	1433	A
1	1	1437	C
1	1	1451	C
1	1	1452	G
1	1	1454	C
1	1	1455	G
1	1	1458	U
1	1	1460	U
1	1	1461	C
1	1	1476	U
1	1	1478	G
1	1	1482	G
1	1	1490	A
1	1	1491	G
1	1	1493	C

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Mol	Chain	Res	Type
1	1	1497	U
1	1	1502	A
1	1	1503	A
1	1	1508	A
1	1	1509	A
1	1	1510	G
1	1	1515	A
1	1	1524	G
1	1	1528	A
1	1	1529	G
1	1	1530	G
1	1	1532	A
1	1	1534	U
1	1	1535	A
1	1	1536	C
1	1	1538	G
1	1	1542	U
1	1	1544	A
1	1	1554	U
1	1	1558	C
1	1	1560	G
1	1	1566	A
1	1	1569	A
1	1	1578	U
1	1	1580	A
1	1	1581	G
1	1	1582	C
1	1	1583	A
1	1	1584	U
1	1	1585	C
1	1	1586	A
1	1	1587	G
1	1	1588	G
1	1	1589	U
1	1	1590	A
1	1	1597	A
1	1	1607	C
1	1	1610	A
1	1	1634	A
1	1	1646	C
1	1	1647	U
1	1	1648	U

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Mol	Chain	Res	Type
1	1	1651	G
1	1	1665	A
1	1	1669	A
1	1	1674	G
1	1	1695	G
1	1	1715	G
1	1	1716	U
1	1	1722	A
1	1	1723	G
1	1	1724	G
1	1	1725	U
1	1	1729	U
1	1	1731	G
1	1	1732	C
1	1	1733	G
1	1	1738	G
1	1	1756	G
1	1	1757	A
1	1	1758	U
1	1	1759	A
1	1	1764	C
1	1	1773	A
1	1	1780	A
1	1	1781	U
1	1	1784	A
1	1	1791	A
1	1	1792	G
1	1	1797	G
1	1	1800	C
1	1	1801	A
1	1	1802	A
1	1	1807	G
1	1	1808	A
1	1	1809	A
1	1	1811	G
1	1	1815	A
1	1	1816	C
1	1	1829	A
1	1	1833	C
1	1	1847	A
1	1	1848	A
1	1	1850	G

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Mol	Chain	Res	Type
1	1	1858	A
1	1	1859	U
1	1	1864	U
1	1	1865	U
1	1	1866	A
1	1	1869	G
1	1	1870	C
1	1	1871	A
1	1	1872	A
1	1	1873	G
1	1	1875	G
1	1	1876	A
1	1	1884	G
1	1	1885	A
1	1	1900	A
1	1	1906	G
1	1	1913	A
1	1	1914	C
1	1	1923	U
1	1	1924	C
1	1	1926	U
1	1	1929	G
1	1	1930	G
1	1	1936	A
1	1	1940	U
1	1	1955	U
1	1	1963	U
1	1	1967	C
1	1	1970	A
1	1	1971	U
1	1	1972	G
1	1	1982	U
1	1	1991	U
1	1	1992	G
1	1	1993	U
1	1	1997	C
1	1	2002	G
1	1	2020	A
1	1	2022	U
1	1	2023	C
1	1	2030	6MZ
1	1	2031	A

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Mol	Chain	Res	Type
1	1	2033	A
1	1	2043	C
1	1	2055	C
1	1	2056	G
1	1	2060	A
1	1	2061	G
1	1	2062	A
1	1	2069	G7M
1	1	2072	C
1	1	2077	A
1	1	2093	G
1	1	2099	U
1	1	2101	A
1	1	2102	G
1	1	2110	G
1	1	2111	U
1	1	2112	G
1	1	2113	U
1	1	2114	A
1	1	2115	G
1	1	2116	G
1	1	2117	A
1	1	2119	A
1	1	2120	G
1	1	2122	U
1	1	2125	G
1	1	2126	A
1	1	2127	G
1	1	2129	C
1	1	2131	U
1	1	2132	U
1	1	2133	G
1	1	2134	A
1	1	2137	U
1	1	2138	G
1	1	2139	U
1	1	2140	G
1	1	2142	A
1	1	2145	C
1	1	2146	C
1	1	2147	A
1	1	2153	C

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Mol	Chain	Res	Type
1	1	2154	A
1	1	2158	A
1	1	2163	A
1	1	2164	C
1	1	2165	C
1	1	2167	U
1	1	2170	A
1	1	2171	A
1	1	2172	U
1	1	2173	A
1	1	2176	A
1	1	2178	C
1	1	2181	U
1	1	2182	U
1	1	2183	A
1	1	2188	U
1	1	2189	U
1	1	2190	G
1	1	2193	G
1	1	2194	U
1	1	2198	A
1	1	2199	A
1	1	2204	G
1	1	2211	A
1	1	2212	A
1	1	2225	A
1	1	2226	C
1	1	2238	G
1	1	2239	G
1	1	2243	U
1	1	2250	G
1	1	2266	A
1	1	2269	G
1	1	2278	A
1	1	2279	G
1	1	2283	C
1	1	2287	A
1	1	2288	A
1	1	2296	U
1	1	2305	U
1	1	2309	A
1	1	2315	G

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Mol	Chain	Res	Type
1	1	2319	G
1	1	2320	U
1	1	2325	G
1	1	2327	A
1	1	2331	G
1	1	2333	A
1	1	2335	A
1	1	2336	A
1	1	2343	U
1	1	2344	U
1	1	2347	C
1	1	2350	C
1	1	2357	G
1	1	2361	G
1	1	2376	A
1	1	2383	G
1	1	2385	C
1	1	2388	A
1	1	2391	G
1	1	2402	U
1	1	2403	C
1	1	2406	A
1	1	2422	C
1	1	2423	U
1	1	2424	C
1	1	2425	A
1	1	2426	A
1	1	2428	G
1	1	2429	G
1	1	2430	A
1	1	2431	U
1	1	2441	U
1	1	2445	2MG
1	1	2447	G
1	1	2448	A
1	1	2459	A
1	1	2469	A
1	1	2470	G
1	1	2475	C
1	1	2476	A
1	1	2478	A
1	1	2484	G

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Mol	Chain	Res	Type
1	1	2490	G
1	1	2491	U
1	1	2492	U
1	1	2494	G
1	1	2501	C
1	1	2502	G
1	1	2503	2MA
1	1	2504	PSU
1	1	2507	C
1	1	2513	A
1	1	2518	A
1	1	2520	C
1	1	2529	G
1	1	2534	A
1	1	2535	G
1	1	2543	G
1	1	2547	A
1	1	2548	U
1	1	2554	U
1	1	2558	C
1	1	2566	A
1	1	2567	G
1	1	2572	A
1	1	2573	C
1	1	2574	G
1	1	2585	U
1	1	2586	U
1	1	2602	A
1	1	2608	G
1	1	2609	U
1	1	2610	C
1	1	2611	C
1	1	2613	U
1	1	2614	A
1	1	2615	U
1	1	2621	G
1	1	2629	U
1	1	2640	G
1	1	2642	G
1	1	2689	U
1	1	2690	U
1	1	2714	G

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Mol	Chain	Res	Type
1	1	2720	U
1	1	2725	A
1	1	2726	A
1	1	2728	U
1	1	2732	G
1	1	2733	A
1	1	2739	U
1	1	2744	G
1	1	2748	A
1	1	2757	A
1	1	2765	A
1	1	2766	A
1	1	2776	A
1	1	2778	A
1	1	2780	G
1	1	2793	C
1	1	2794	C
1	1	2797	U
1	1	2798	U
1	1	2799	A
1	1	2800	A
1	1	2811	G
1	1	2818	U
1	1	2820	A
1	1	2821	A
1	1	2825	G
1	1	2833	U
1	1	2835	A
1	1	2836	U
1	1	2861	U
1	1	2867	G
1	1	2872	A
1	1	2873	A
1	1	2879	A
1	1	2880	C
1	1	2883	A
1	1	2884	U
1	1	2885	G
1	1	2893	A
1	1	2901	C
1	1	2902	C
2	2	3	A

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Mol	Chain	Res	Type
2	2	4	U
2	2	5	U
2	2	6	G
2	2	7	A
2	2	8	A
2	2	9	G
2	2	22	G
2	2	32	A
2	2	33	A
2	2	39	G
2	2	44	A
2	2	47	C
2	2	48	C
2	2	50	A
2	2	51	A
2	2	52	C
2	2	54	C
2	2	65	A
2	2	68	G
2	2	70	U
2	2	71	A
2	2	72	A
2	2	74	A
2	2	77	A
2	2	79	G
2	2	81	A
2	2	82	G
2	2	83	C
2	2	84	U
2	2	85	U
2	2	86	G
2	2	87	C
2	2	88	U
2	2	89	U
2	2	91	U
2	2	92	U
2	2	94	G
2	2	95	C
2	2	108	G
2	2	116	A
2	2	120	A
2	2	121	U

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Mol	Chain	Res	Type
2	2	129	A
2	2	130	A
2	2	131	A
2	2	141	G
2	2	142	G
2	2	149	A
2	2	163	C
2	2	164	G
2	2	168	G
2	2	173	U
2	2	177	G
2	2	181	A
2	2	182	A
2	2	183	C
2	2	185	U
2	2	189	A
2	2	192	A
2	2	197	A
2	2	202	G
2	2	204	G
2	2	209	U
2	2	211	G
2	2	212	G
2	2	213	G
2	2	214	C
2	2	226	G
2	2	230	G
2	2	244	U
2	2	245	U
2	2	247	G
2	2	251	G
2	2	253	A
2	2	266	G
2	2	267	C
2	2	276	G
2	2	279	A
2	2	289	G
2	2	301	G
2	2	306	A
2	2	319	G
2	2	328	C
2	2	329	A

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Mol	Chain	Res	Type
2	2	330	C
2	2	332	G
2	2	338	A
2	2	347	G
2	2	348	G
2	2	351	G
2	2	352	C
2	2	353	A
2	2	354	G
2	2	356	A
2	2	366	A
2	2	367	U
2	2	372	C
2	2	382	A
2	2	397	A
2	2	406	G
2	2	409	U
2	2	411	A
2	2	412	A
2	2	413	G
2	2	414	A
2	2	415	A
2	2	421	U
2	2	422	C
2	2	423	G
2	2	424	G
2	2	428	G
2	2	429	U
2	2	431	A
2	2	438	U
2	2	448	A
2	2	452	A
2	2	453	G
2	2	456	A
2	2	457	G
2	2	458	U
2	2	464	U
2	2	467	U
2	2	468	A
2	2	476	U
2	2	477	C
2	2	478	A

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Mol	Chain	Res	Type
2	2	479	U
2	2	481	G
2	2	486	U
2	2	495	A
2	2	496	A
2	2	497	G
2	2	509	A
2	2	510	A
2	2	511	C
2	2	515	G
2	2	517	G
2	2	518	C
2	2	519	C
2	2	521	G
2	2	527	7MG
2	2	531	U
2	2	532	A
2	2	533	A
2	2	534	U
2	2	547	A
2	2	562	U
2	2	572	A
2	2	573	A
2	2	576	C
2	2	577	G
2	2	585	G
2	2	596	A
2	2	607	A
2	2	615	G
2	2	616	G
2	2	618	C
2	2	619	U
2	2	633	G
2	2	639	G
2	2	640	A
2	2	641	U
2	2	652	U
2	2	653	U
2	2	661	G
2	2	665	A
2	2	682	G
2	2	686	U

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Mol	Chain	Res	Type
2	2	695	A
2	2	701	U
2	2	703	G
2	2	718	A
2	2	723	U
2	2	724	G
2	2	731	G
2	2	734	G
2	2	744	C
2	2	747	A
2	2	748	G
2	2	755	G
2	2	793	U
2	2	794	A
2	2	799	G
2	2	809	G
2	2	814	A
2	2	815	A
2	2	817	C
2	2	819	A
2	2	821	G
2	2	828	U
2	2	829	G
2	2	832	G
2	2	842	U
2	2	843	U
2	2	844	G
2	2	846	G
2	2	849	G
2	2	864	A
2	2	871	U
2	2	889	A
2	2	902	G
2	2	914	A
2	2	926	G
2	2	934	C
2	2	935	A
2	2	958	A
2	2	960	U
2	2	966	2MG
2	2	967	5MC
2	2	968	A

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Mol	Chain	Res	Type
2	2	969	A
2	2	971	G
2	2	974	A
2	2	975	A
2	2	976	G
2	2	977	A
2	2	982	U
2	2	987	G
2	2	989	U
2	2	993	G
2	2	996	A
2	2	999	C
2	2	1000	A
2	2	1003	G
2	2	1004	A
2	2	1008	U
2	2	1009	U
2	2	1017	U
2	2	1018	G
2	2	1019	A
2	2	1020	G
2	2	1022	A
2	2	1023	U
2	2	1024	G
2	2	1026	G
2	2	1028	C
2	2	1029	U
2	2	1032	G
2	2	1033	G
2	2	1034	G
2	2	1036	A
2	2	1043	G
2	2	1044	A
2	2	1045	C
2	2	1046	A
2	2	1054	C
2	2	1055	A
2	2	1065	U
2	2	1067	A
2	2	1070	U
2	2	1092	A
2	2	1094	G

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Mol	Chain	Res	Type
2	2	1095	U
2	2	1099	G
2	2	1101	A
2	2	1102	A
2	2	1108	G
2	2	1124	G
2	2	1125	U
2	2	1126	U
2	2	1132	C
2	2	1133	G
2	2	1136	C
2	2	1137	C
2	2	1139	G
2	2	1140	C
2	2	1141	C
2	2	1145	A
2	2	1146	A
2	2	1151	A
2	2	1152	A
2	2	1159	U
2	2	1160	G
2	2	1167	A
2	2	1168	U
2	2	1169	A
2	2	1170	A
2	2	1171	A
2	2	1176	A
2	2	1177	G
2	2	1184	G
2	2	1193	G
2	2	1196	A
2	2	1197	A
2	2	1201	A
2	2	1212	U
2	2	1213	A
2	2	1214	C
2	2	1225	A
2	2	1226	C
2	2	1227	A
2	2	1228	C
2	2	1233	G
2	2	1238	A

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Mol	Chain	Res	Type
2	2	1240	U
2	2	1241	G
2	2	1256	A
2	2	1257	A
2	2	1258	G
2	2	1260	G
2	2	1261	A
2	2	1275	A
2	2	1279	G
2	2	1280	A
2	2	1285	A
2	2	1287	A
2	2	1297	G
2	2	1299	A
2	2	1300	G
2	2	1301	U
2	2	1302	C
2	2	1305	G
2	2	1312	G
2	2	1317	C
2	2	1319	A
2	2	1320	C
2	2	1323	G
2	2	1331	G
2	2	1336	C
2	2	1340	A
2	2	1348	U
2	2	1353	G
2	2	1363	A
2	2	1368	A
2	2	1370	G
2	2	1378	C
2	2	1379	G
2	2	1381	U
2	2	1397	C
2	2	1398	A
2	2	1399	C
2	2	1401	G
2	2	1406	U
2	2	1419	G
2	2	1429	A
2	2	1432	G

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Mol	Chain	Res	Type
2	2	1433	A
2	2	1441	A
2	2	1445	U
2	2	1446	A
2	2	1451	U
2	2	1452	C
2	2	1453	G
2	2	1462	C
2	2	1471	U
2	2	1484	C
2	2	1487	G
2	2	1491	G
2	2	1492	A
2	2	1493	A
2	2	1494	G
2	2	1497	G
2	2	1499	A
2	2	1503	A
2	2	1506	U
2	2	1517	G
2	2	1519	MA6
2	2	1520	C
2	2	1529	G
2	2	1530	G
2	2	1533	C
2	2	1534	A
3	3	2	G
3	3	12	C
3	3	13	G
3	3	15	A
3	3	16	G
3	3	17	C
3	3	19	C
3	3	25	U
3	3	35	C
3	3	36	C
3	3	40	U
3	3	41	G
3	3	44	G
3	3	51	G
3	3	52	A
3	3	53	A

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Mol	Chain	Res	Type
3	3	54	G
3	3	56	G
3	3	58	A
3	3	66	A
3	3	67	G
3	3	68	C
3	3	88	C
3	3	89	U
3	3	90	C
3	3	91	C
3	3	99	A
3	3	105	G
3	3	108	A
3	3	109	A
52	5	3	G
52	5	7	U
52	5	8	4SU
52	5	9	G
52	5	12	C
52	5	15	C
52	5	16	C
52	5	17	U
52	5	18	G
52	5	19	G
52	5	20	H2U
52	5	21	A
52	5	22	G
52	5	47	U
52	5	48	C
52	5	49	G
52	5	52	G
52	5	55	PSU
52	5	58	A
52	5	59	A
52	5	68	C
52	5	70	G
52	5	72	A
52	5	76	A
55	4	19	U
55	4	21	A

All (22) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	1	404	A
1	1	479	A
1	1	545	U
1	1	546	U
1	1	613	A
1	1	647	G
1	1	654	A
1	1	784	G
1	1	830	G
1	1	894	U
1	1	1379	U
1	1	2146	C
1	1	2193	G
1	1	2425	A
1	1	2756	U
2	2	496	A
2	2	516	PSU
2	2	966	2MG
2	2	1109	C
2	2	1136	C
2	2	1145	A
52	5	17	U

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

40 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
2	4OC	2	1402	2	20,23,24	2.87	8 (40%)	25,32,35	1.15	3 (12%)
2	5MC	2	967	2	19,22,23	3.83	8 (42%)	26,32,35	1.12	3 (11%)
1	PSU	1	1917	1	18,21,22	1.09	1 (5%)	21,30,33	2.03	4 (19%)
1	G7M	1	2069	1	23,26,27	2.52	7 (30%)	34,39,42	2.43	10 (29%)
52	4OC	5	32	52	20,23,24	3.02	8 (40%)	25,32,35	1.01	2 (8%)
1	2MG	1	1835	1	23,26,27	2.89	8 (34%)	33,38,41	2.36	9 (27%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	6MZ	1	2030	1	22,25,26	2.85	7 (31%)	29,36,39	4.61	14 (48%)
2	2MG	2	966	2	23,26,27	2.97	7 (30%)	33,38,41	2.42	7 (21%)
2	MA6	2	1518	2	23,26,27	1.63	5 (21%)	33,38,41	2.73	12 (36%)
2	MA6	2	1519	2	23,26,27	1.55	4 (17%)	33,38,41	2.74	13 (39%)
1	PSU	1	746	1,56	18,21,22	1.12	2 (11%)	21,30,33	2.09	6 (28%)
1	2MG	1	2445	1	23,26,27	2.83	7 (30%)	33,38,41	2.42	9 (27%)
42	0TD	q	89	42	8,9,10	1.83	2 (25%)	6,11,13	2.25	2 (33%)
1	OMG	1	2251	52,1	23,26,27	2.35	7 (30%)	32,38,41	1.95	8 (25%)
2	7MG	2	527	2	23,26,27	3.29	10 (43%)	27,39,42	2.00	8 (29%)
2	5MC	2	1407	2	19,22,23	3.69	8 (42%)	26,32,35	1.23	3 (11%)
52	4SU	5	8	52	18,21,22	3.69	7 (38%)	25,30,33	2.20	5 (20%)
52	H2U	5	20	52	18,21,22	3.31	5 (27%)	19,30,33	1.35	3 (15%)
1	1MG	1	745	1	23,26,27	2.75	6 (26%)	33,39,42	2.25	9 (27%)
1	PSU	1	2605	1	18,21,22	1.09	2 (11%)	21,30,33	2.16	6 (28%)
1	PSU	1	2457	1	18,21,22	1.18	3 (16%)	21,30,33	2.44	6 (28%)
53	FME	H	77	53	8,9,10	0.70	0	8,9,11	1.28	1 (12%)
1	5MU	1	747	1	19,22,23	4.48	7 (36%)	27,32,35	3.92	12 (44%)
1	PSU	1	2580	1	18,21,22	1.20	3 (16%)	21,30,33	2.04	5 (23%)
1	PSU	1	955	1	18,21,22	1.09	2 (11%)	21,30,33	1.82	4 (19%)
1	5MC	1	1962	1	19,22,23	3.62	8 (42%)	26,32,35	1.34	3 (11%)
1	6MZ	1	1618	1	22,25,26	2.86	6 (27%)	29,36,39	4.47	12 (41%)
2	2MG	2	1516	2	23,26,27	2.92	8 (34%)	33,38,41	2.18	11 (33%)
52	5MU	5	54	52	19,22,23	4.81	7 (36%)	27,32,35	3.43	9 (33%)
1	3TD	1	1915	1	19,22,23	4.26	5 (26%)	23,32,35	2.01	3 (13%)
52	PSU	5	55	52	18,21,22	1.20	1 (5%)	21,30,33	1.64	4 (19%)
1	5MU	1	1939	1	19,22,23	4.56	7 (36%)	27,32,35	4.00	10 (37%)
1	PSU	1	1911	1	18,21,22	1.09	2 (11%)	21,30,33	1.96	4 (19%)
2	2MG	2	1207	2	23,26,27	2.84	9 (39%)	33,38,41	2.57	12 (36%)
1	2MA	1	2503	1,56	22,25,26	3.45	10 (45%)	32,37,40	1.96	9 (28%)
1	OMU	1	2552	1	19,22,23	2.80	8 (42%)	25,31,34	1.97	5 (20%)
2	UR3	2	1498	2	19,22,23	2.46	6 (31%)	26,32,35	1.64	3 (11%)
2	PSU	2	516	2	18,21,22	1.05	1 (5%)	21,30,33	1.90	4 (19%)
1	PSU	1	2504	1	18,21,22	1.07	2 (11%)	21,30,33	1.96	4 (19%)
1	OMC	1	2498	1,56	19,22,23	2.76	7 (36%)	25,31,34	0.78	0

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
2	4OC	2	1402	2	-	2/9/29/30	0/2/2/2
2	5MC	2	967	2	-	0/7/25/26	0/2/2/2
1	PSU	1	1917	1	-	0/7/25/26	0/2/2/2
1	G7M	1	2069	1	-	0/7/25/26	0/3/3/3
52	4OC	5	32	52	-	0/9/29/30	0/2/2/2
1	2MG	1	1835	1	-	0/9/27/28	0/3/3/3
1	6MZ	1	2030	1	-	2/9/27/28	0/3/3/3
2	2MG	2	966	2	-	3/9/27/28	0/3/3/3
2	MA6	2	1518	2	-	0/11/29/30	0/3/3/3
2	MA6	2	1519	2	-	3/11/29/30	0/3/3/3
1	PSU	1	746	1,56	-	4/7/25/26	0/2/2/2
1	2MG	1	2445	1	-	2/9/27/28	0/3/3/3
42	0TD	q	89	42	-	1/7/12/14	-
1	OMG	1	2251	52,1	-	0/9/27/28	0/3/3/3
2	7MG	2	527	2	-	3/7/37/38	0/3/3/3
2	5MC	2	1407	2	-	0/7/25/26	0/2/2/2
52	4SU	5	8	52	-	4/7/25/26	0/2/2/2
52	H2U	5	20	52	-	4/7/38/39	0/2/2/2
1	1MG	1	745	1	-	0/7/25/26	0/3/3/3
1	PSU	1	2605	1	-	0/7/25/26	0/2/2/2
1	PSU	1	2457	1	-	2/7/25/26	0/2/2/2
53	FME	H	77	53	-	2/7/9/11	-
1	5MU	1	747	1	-	1/7/25/26	0/2/2/2
1	PSU	1	2580	1	-	0/7/25/26	0/2/2/2
1	PSU	1	955	1	-	0/7/25/26	0/2/2/2
1	5MC	1	1962	1	-	3/7/25/26	0/2/2/2
1	6MZ	1	1618	1	-	2/9/27/28	0/3/3/3
2	2MG	2	1516	2	-	0/9/27/28	0/3/3/3
52	5MU	5	54	52	-	0/7/25/26	0/2/2/2
1	3TD	1	1915	1	-	3/7/25/26	0/2/2/2
52	PSU	5	55	52	-	2/7/25/26	0/2/2/2
1	5MU	1	1939	1	-	0/7/25/26	0/2/2/2
1	PSU	1	1911	1	-	0/7/25/26	0/2/2/2
2	2MG	2	1207	2	-	0/9/27/28	0/3/3/3
1	2MA	1	2503	1,56	-	3/7/25/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OMU	1	2552	1	-	0/9/27/28	0/2/2/2
2	UR3	2	1498	2	-	0/7/25/26	0/2/2/2
2	PSU	2	516	2	-	2/7/25/26	0/2/2/2
1	PSU	1	2504	1	-	1/7/25/26	0/2/2/2
1	OMC	1	2498	1,56	-	0/9/27/28	0/2/2/2

All (221) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	1915	3TD	C6-C5	13.45	1.50	1.35
52	5	54	5MU	C2-N1	11.33	1.56	1.38
1	1	1618	6MZ	C6-N6	10.72	1.46	1.34
52	5	20	H2U	C2-N1	10.62	1.50	1.35
52	5	54	5MU	C6-N1	10.43	1.55	1.38
1	1	2030	6MZ	C6-N6	10.22	1.45	1.34
1	1	747	5MU	C2-N1	9.99	1.54	1.38
1	1	2503	2MA	C4-N3	9.87	1.47	1.34
52	5	54	5MU	C4-C5	9.78	1.60	1.44
1	1	1939	5MU	C2-N1	9.75	1.53	1.38
1	1	1915	3TD	C2-N1	9.49	1.48	1.37
1	1	1939	5MU	C4-C5	9.44	1.60	1.44
1	1	1939	5MU	C6-N1	9.40	1.54	1.38
1	1	747	5MU	C6-N1	9.32	1.53	1.38
2	2	1407	5MC	C6-C5	9.09	1.49	1.34
2	2	967	5MC	C6-C5	8.92	1.49	1.34
1	1	747	5MU	C4-C5	8.74	1.59	1.44
1	1	747	5MU	C4-N3	-8.52	1.22	1.38
1	1	1962	5MC	C6-C5	8.46	1.48	1.34
1	1	1939	5MU	C4-N3	-8.45	1.23	1.38
52	5	8	4SU	C4-N3	7.96	1.45	1.37
52	5	8	4SU	C2-N1	7.91	1.50	1.38
52	5	54	5MU	C4-N3	-7.83	1.24	1.38
2	2	966	2MG	C2-N3	7.47	1.46	1.32
1	1	745	1MG	C2-N2	7.37	1.47	1.34
1	1	1962	5MC	C4-N3	7.27	1.45	1.34
2	2	1516	2MG	C2-N3	7.19	1.45	1.32
1	1	1835	2MG	C2-N3	7.11	1.45	1.32
2	2	527	7MG	C8-N9	7.08	1.50	1.45
2	2	967	5MC	C4-N3	7.07	1.45	1.34
52	5	32	4OC	C4-N3	6.95	1.44	1.32
1	1	2445	2MG	C2-N3	6.94	1.45	1.32
2	2	966	2MG	C4-N3	6.92	1.50	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2	527	7MG	C5-N7	6.81	1.44	1.35
52	5	20	H2U	C2-N3	6.70	1.49	1.38
2	2	966	2MG	C2-N2	6.61	1.47	1.33
2	2	1207	2MG	C2-N3	6.59	1.44	1.32
1	1	1835	2MG	C4-N3	6.55	1.49	1.34
2	2	1402	4OC	C4-N3	6.54	1.43	1.32
2	2	1516	2MG	C2-N2	6.45	1.46	1.33
1	1	2552	OMU	C2-N3	6.36	1.49	1.38
2	2	1407	5MC	C4-N3	6.28	1.44	1.34
1	1	745	1MG	C2-N3	6.27	1.43	1.33
2	2	1516	2MG	C4-N3	6.25	1.48	1.34
1	1	1835	2MG	C2-N2	6.24	1.46	1.33
1	1	2445	2MG	C4-N3	6.24	1.48	1.34
2	2	1207	2MG	C2-N2	6.23	1.46	1.33
1	1	1962	5MC	C2-N3	6.23	1.48	1.36
1	1	2503	2MA	C2-N3	6.13	1.44	1.34
2	2	967	5MC	C2-N3	6.11	1.48	1.36
2	2	967	5MC	C5-C4	6.10	1.48	1.44
52	5	32	4OC	C6-C5	6.05	1.49	1.35
1	1	2498	OMC	C6-C5	5.97	1.48	1.35
1	1	2445	2MG	C2-N2	5.96	1.45	1.33
1	1	2069	G7M	C4-N3	5.94	1.47	1.34
1	1	2552	OMU	C2-N1	5.94	1.47	1.38
1	1	745	1MG	C4-N3	5.92	1.47	1.34
2	2	1498	UR3	C2-N1	5.92	1.46	1.38
1	1	2069	G7M	C5-N7	-5.91	1.32	1.39
2	2	1207	2MG	C4-N3	5.90	1.47	1.34
52	5	8	4SU	C6-C5	5.86	1.48	1.35
1	1	2251	OMG	C4-N3	5.83	1.47	1.34
52	5	32	4OC	C2-N3	5.81	1.47	1.36
1	1	2498	OMC	C2-N3	5.76	1.47	1.36
2	2	1498	UR3	C6-C5	5.72	1.48	1.35
2	2	1407	5MC	C2-N3	5.71	1.47	1.36
2	2	1402	4OC	C6-C5	5.70	1.48	1.35
52	5	8	4SU	C2-N3	5.65	1.47	1.38
52	5	8	4SU	C4-S4	-5.64	1.58	1.68
2	2	527	7MG	C2-N3	5.64	1.46	1.33
1	1	1915	3TD	C6-N1	5.62	1.45	1.36
2	2	1402	4OC	C2-N3	5.50	1.47	1.36
52	5	54	5MU	C6-C5	5.48	1.43	1.34
1	1	2503	2MA	C2-N1	5.47	1.43	1.34
2	2	1407	5MC	C5-C4	5.47	1.48	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	2552	OMU	C6-C5	5.45	1.47	1.35
1	1	1939	5MU	C6-C5	5.45	1.43	1.34
2	2	527	7MG	C4-N3	5.40	1.46	1.34
1	1	2498	OMC	C4-N3	5.39	1.45	1.34
1	1	2503	2MA	C6-N1	5.25	1.42	1.35
52	5	20	H2U	C4-N3	5.04	1.46	1.37
1	1	1915	3TD	C2-N3	4.99	1.49	1.38
1	1	2251	OMG	C2-N3	4.93	1.45	1.33
1	1	747	5MU	C6-C5	4.88	1.42	1.34
2	2	967	5MC	C4-N4	4.82	1.46	1.34
2	2	527	7MG	C2-N2	4.75	1.45	1.34
2	2	1407	5MC	C4-N4	4.74	1.46	1.34
1	1	2069	G7M	C2-N3	4.69	1.44	1.33
2	2	1207	2MG	C2-N1	4.61	1.44	1.36
1	1	1962	5MC	C4-N4	4.61	1.45	1.34
2	2	527	7MG	C4-N9	4.53	1.43	1.37
1	1	2251	OMG	C2-N2	4.49	1.44	1.34
1	1	1962	5MC	C5-C4	4.47	1.47	1.44
2	2	1407	5MC	C6-N1	4.41	1.45	1.38
2	2	1498	UR3	C2-N3	4.39	1.47	1.39
2	2	967	5MC	C2-N1	4.39	1.49	1.40
1	1	2030	6MZ	C5-C4	-4.31	1.31	1.39
2	2	1516	2MG	C2-N1	4.25	1.43	1.36
2	2	1407	5MC	C2-N1	4.21	1.48	1.40
2	2	967	5MC	C6-N1	4.18	1.45	1.38
52	5	55	PSU	C6-C5	4.16	1.39	1.35
52	5	32	4OC	C2-N1	4.15	1.48	1.40
2	2	1402	4OC	C2-N1	4.12	1.48	1.40
1	1	2445	2MG	O6-C6	-4.06	1.15	1.23
2	2	966	2MG	C2-N1	4.06	1.43	1.36
1	1	2445	2MG	C5-N7	-4.03	1.31	1.39
1	1	2069	G7M	C2-N2	3.96	1.43	1.34
52	5	32	4OC	C4-N4	3.95	1.44	1.36
1	1	1618	6MZ	C5-C4	-3.95	1.32	1.39
1	1	1962	5MC	C2-N1	3.91	1.48	1.40
1	1	1962	5MC	C6-N1	3.91	1.44	1.38
1	1	2030	6MZ	C8-N9	-3.91	1.30	1.37
1	1	1835	2MG	C2-N1	3.78	1.42	1.36
1	1	1835	2MG	C5-N7	-3.74	1.31	1.39
1	1	2503	2MA	C5-N7	-3.73	1.32	1.39
1	1	2503	2MA	C5-C6	3.71	1.51	1.41
1	1	2503	2MA	C6-N6	-3.70	1.24	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	2552	OMU	C4-N3	3.67	1.44	1.38
2	2	1519	MA6	C6-N6	3.64	1.46	1.36
1	1	1835	2MG	O6-C6	-3.64	1.16	1.23
2	2	1402	4OC	C4-N4	3.63	1.43	1.36
1	1	1618	6MZ	C8-N9	-3.63	1.31	1.37
1	1	2030	6MZ	C4-N9	-3.60	1.30	1.37
2	2	1516	2MG	C5-N7	-3.60	1.31	1.39
1	1	2251	OMG	C5-N7	-3.59	1.31	1.39
52	5	32	4OC	C5-C4	3.57	1.48	1.41
2	2	966	2MG	C5-N7	-3.57	1.31	1.39
2	2	1518	MA6	C8-N9	-3.57	1.31	1.37
2	2	1518	MA6	C6-N6	3.55	1.46	1.36
2	2	1516	2MG	O6-C6	-3.55	1.16	1.23
2	2	527	7MG	C5-C6	3.52	1.52	1.43
2	2	527	7MG	C2-N1	3.47	1.46	1.37
1	1	2498	OMC	C4-N4	3.47	1.42	1.33
2	2	1207	2MG	O6-C6	-3.45	1.17	1.23
1	1	2069	G7M	C5-C6	3.45	1.52	1.43
1	1	2503	2MA	C4-N9	-3.43	1.30	1.37
2	2	966	2MG	O6-C6	-3.42	1.17	1.23
1	1	745	1MG	C5-N7	-3.41	1.32	1.39
1	1	1618	6MZ	C4-N9	-3.38	1.30	1.37
2	2	1207	2MG	C5-N7	-3.38	1.32	1.39
2	2	1407	5MC	O2-C2	-3.36	1.17	1.23
1	1	1962	5MC	O2-C2	-3.30	1.17	1.23
2	2	1518	MA6	C5-C4	-3.30	1.33	1.39
52	5	8	4SU	C5-C4	3.30	1.46	1.42
1	1	2498	OMC	C6-N1	3.28	1.45	1.38
2	2	967	5MC	O2-C2	-3.27	1.17	1.23
1	1	1911	PSU	C6-C5	3.24	1.38	1.35
2	2	1519	MA6	C5-N7	-3.22	1.33	1.39
2	2	1402	4OC	C5-C4	3.22	1.48	1.41
1	1	2552	OMU	O4-C4	-3.22	1.18	1.24
2	2	1518	MA6	C5-N7	-3.21	1.33	1.39
1	1	2498	OMC	O2-C2	-3.19	1.17	1.23
1	1	1939	5MU	O4-C4	-3.17	1.17	1.23
1	1	2445	2MG	C2-N1	3.14	1.41	1.36
1	1	2498	OMC	C2-N1	3.14	1.46	1.40
42	q	89	0TD	CB-CA	-3.12	1.53	1.54
1	1	1917	PSU	C6-C5	3.12	1.38	1.35
1	1	747	5MU	O2-C2	-3.10	1.17	1.23
2	2	516	PSU	C6-C5	3.09	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	2069	G7M	O6-C6	-3.07	1.17	1.23
2	2	1519	MA6	C8-N9	-3.06	1.32	1.37
52	5	32	4OC	C6-N1	3.03	1.45	1.38
1	1	745	1MG	C2-N1	3.03	1.42	1.37
1	1	746	PSU	C6-C5	3.02	1.38	1.35
1	1	747	5MU	O4-C4	-3.02	1.17	1.23
1	1	2030	6MZ	C6-N1	-2.98	1.30	1.35
1	1	2251	OMG	O6-C6	-2.97	1.17	1.23
1	1	1939	5MU	O2-C2	-2.97	1.17	1.23
2	2	1519	MA6	C5-C4	-2.95	1.33	1.39
1	1	1915	3TD	C4-N3	2.93	1.46	1.40
2	2	1402	4OC	O2-C2	-2.90	1.18	1.23
1	1	1618	6MZ	C6-N1	-2.86	1.30	1.35
2	2	527	7MG	C6-N1	2.80	1.44	1.38
52	5	32	4OC	O2-C2	-2.79	1.18	1.23
2	2	527	7MG	O6-C6	-2.73	1.18	1.23
2	2	1402	4OC	C6-N1	2.70	1.44	1.38
52	5	54	5MU	O4-C4	-2.70	1.18	1.23
1	1	2580	PSU	C6-C5	2.68	1.38	1.35
1	1	2552	OMU	O2-C2	-2.67	1.18	1.23
1	1	2030	6MZ	C5-C6	-2.65	1.35	1.41
1	1	2605	PSU	C6-C5	2.64	1.38	1.35
1	1	2504	PSU	C6-C5	2.63	1.38	1.35
1	1	955	PSU	C6-C5	2.61	1.38	1.35
1	1	745	1MG	C4-N9	-2.54	1.31	1.38
1	1	2457	PSU	C6-C5	2.54	1.38	1.35
52	5	8	4SU	O2-C2	-2.53	1.18	1.23
1	1	2503	2MA	C8-N9	-2.52	1.33	1.37
1	1	2580	PSU	C4-C5	-2.52	1.37	1.44
1	1	2580	PSU	O4'-C1'	-2.51	1.40	1.43
2	2	1498	UR3	O4-C4	-2.51	1.18	1.23
2	2	1498	UR3	O2-C2	-2.48	1.17	1.22
1	1	2251	OMG	C4-N9	-2.48	1.31	1.38
1	1	1618	6MZ	C5-C6	-2.47	1.35	1.41
1	1	2457	PSU	C4-C5	-2.45	1.37	1.44
2	2	1207	2MG	C6-N1	2.44	1.43	1.38
1	1	2457	PSU	O4'-C1'	-2.42	1.40	1.43
52	5	54	5MU	O2-C2	-2.42	1.18	1.23
1	1	2504	PSU	C4-C5	-2.40	1.37	1.44
2	2	1207	2MG	C4-N9	-2.37	1.32	1.38
52	5	20	H2U	O4-C4	-2.30	1.18	1.23
1	1	2552	OMU	C6-N1	2.30	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
52	5	20	H2U	O2-C2	-2.28	1.19	1.23
1	1	2251	OMG	C2-N1	2.27	1.43	1.37
1	1	746	PSU	C4-C5	-2.26	1.38	1.44
2	2	1207	2MG	C5-C6	2.26	1.52	1.44
2	2	1498	UR3	C6-N1	2.25	1.43	1.38
1	1	955	PSU	C4-C5	-2.25	1.38	1.44
1	1	2605	PSU	C4-C5	-2.23	1.38	1.44
2	2	1516	2MG	C4-N9	-2.19	1.32	1.38
2	2	966	2MG	C5-C6	2.19	1.52	1.44
1	1	1835	2MG	C5-C6	2.18	1.52	1.44
42	q	89	0TD	CSB-SB	-2.14	1.75	1.79
1	1	2552	OMU	C5-C4	2.11	1.48	1.43
2	2	1516	2MG	C5-C6	2.09	1.52	1.44
1	1	2030	6MZ	C5-N7	-2.08	1.35	1.39
1	1	1911	PSU	C4-C5	-2.08	1.38	1.44
1	1	2503	2MA	C5-C4	-2.02	1.35	1.39
1	1	1835	2MG	C8-N9	-2.01	1.33	1.37
1	1	2445	2MG	C8-N9	-2.01	1.33	1.37
2	2	1518	MA6	C5-C6	-2.01	1.36	1.41
1	1	2069	G7M	C2-N1	2.01	1.42	1.37

All (257) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	2030	6MZ	C4-N9-C8	14.53	120.99	105.74
1	1	1618	6MZ	C4-N9-C8	13.72	120.14	105.74
1	1	1939	5MU	C5-C4-N3	12.60	126.28	115.32
1	1	747	5MU	C5-C4-N3	12.15	125.88	115.32
52	5	54	5MU	C5-C4-N3	11.65	125.44	115.32
1	1	1939	5MU	C5-C6-N1	-11.25	111.09	123.31
1	1	747	5MU	C5-C6-N1	-10.97	111.39	123.31
1	1	2030	6MZ	N9-C8-N7	-10.90	98.48	113.94
1	1	1618	6MZ	N9-C8-N7	-10.07	99.65	113.94
1	1	1618	6MZ	C5-C6-N1	9.12	127.94	118.15
1	1	2030	6MZ	C5-C6-N1	8.93	127.74	118.15
2	2	1519	MA6	N1-C6-N6	-8.38	106.64	116.86
2	2	1518	MA6	N1-C6-N6	-8.32	106.72	116.86
52	5	54	5MU	C5-C6-N1	-8.11	114.50	123.31
52	5	8	4SU	C4-N3-C2	-7.23	120.39	127.31
2	2	966	2MG	C2-N3-C4	7.16	120.96	112.00
1	1	1835	2MG	C2-N3-C4	6.97	120.72	112.00
2	2	966	2MG	C5-C4-N3	-6.78	117.60	128.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	2445	2MG	C2-N3-C4	6.76	120.46	112.00
2	2	1207	2MG	N1-C2-N2	6.70	123.40	116.56
1	1	1835	2MG	C5-C4-N3	-6.43	118.15	128.39
1	1	1618	6MZ	N3-C4-N9	6.38	138.02	127.17
1	1	2552	OMU	C4-N3-C2	-6.30	118.80	126.61
1	1	2445	2MG	C5-C4-N3	-6.23	118.48	128.39
1	1	747	5MU	O4-C4-C5	-6.22	117.80	124.92
1	1	2069	G7M	CN7-N7-C5	6.17	134.49	126.80
2	2	1519	MA6	N1-C2-N3	-6.14	119.29	128.58
1	1	2457	PSU	N1-C2-N3	6.12	121.63	115.17
2	2	1207	2MG	C2-N3-C4	6.12	119.65	112.00
1	1	2030	6MZ	N3-C4-N9	6.07	137.49	127.17
2	2	1518	MA6	N1-C2-N3	-6.01	119.48	128.58
1	1	1939	5MU	C4-N3-C2	-6.00	119.47	127.34
1	1	1618	6MZ	C9-N6-C6	-5.97	117.32	122.85
2	2	1516	2MG	C2-N3-C4	5.95	119.44	112.00
1	1	2457	PSU	C4-N3-C2	-5.87	118.29	126.37
1	1	2030	6MZ	N1-C2-N3	-5.77	119.85	128.58
1	1	2445	2MG	C2-N1-C6	-5.76	117.58	124.55
1	1	745	1MG	C1'-N9-C4	-5.76	109.48	126.49
1	1	1915	3TD	C1'-C5-C4	5.67	126.22	117.61
1	1	2605	PSU	C4-N3-C2	-5.67	118.56	126.37
1	1	2503	2MA	C5-C4-N3	-5.62	121.26	127.18
2	2	1519	MA6	C5-C6-N6	5.55	134.12	125.33
1	1	746	PSU	C4-N3-C2	-5.55	118.73	126.37
1	1	1618	6MZ	N1-C2-N3	-5.50	120.26	128.58
1	1	747	5MU	C4-N3-C2	-5.47	120.17	127.34
2	2	1498	UR3	C4-N3-C2	-5.43	120.21	124.58
1	1	2605	PSU	N1-C2-N3	5.40	120.86	115.17
1	1	1915	3TD	N1-C2-N3	5.39	120.05	116.13
1	1	1939	5MU	O4-C4-C5	-5.39	118.75	124.92
1	1	1917	PSU	N1-C2-N3	5.38	120.84	115.17
1	1	2030	6MZ	C9-N6-C6	-5.36	117.88	122.85
1	1	2251	OMG	C5-C4-N3	-5.35	119.88	128.39
1	1	1618	6MZ	C5-C4-N9	-5.27	100.07	105.81
2	2	1518	MA6	C5-C6-N6	5.27	133.67	125.33
1	1	2030	6MZ	C5-C4-N9	-5.26	100.08	105.81
52	5	8	4SU	C5-C4-N3	5.23	119.62	114.75
2	2	1516	2MG	C5-C4-N3	-5.21	120.10	128.39
1	1	1835	2MG	C2-N1-C6	-5.20	118.27	124.55
52	5	54	5MU	O4-C4-C5	-5.15	119.03	124.92
1	1	1939	5MU	N3-C2-N1	5.09	121.52	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	746	PSU	N1-C2-N3	5.08	120.52	115.17
1	1	2069	G7M	C2-N3-C4	5.07	121.04	112.30
1	1	745	1MG	C1'-N9-C8	5.06	141.11	126.73
1	1	745	1MG	C5-C4-N3	-5.06	120.34	128.39
1	1	2580	PSU	C4-N3-C2	-5.06	119.41	126.37
1	1	2069	G7M	CN7-N7-C8	-4.99	117.24	124.79
1	1	2580	PSU	N1-C2-N3	4.98	120.42	115.17
2	2	966	2MG	C2-N1-C6	-4.97	118.55	124.55
2	2	966	2MG	N9-C4-N3	4.95	135.85	125.95
2	2	516	PSU	C4-N3-C2	-4.94	119.57	126.37
1	1	1911	PSU	N1-C2-N3	4.94	120.38	115.17
1	1	1911	PSU	C4-N3-C2	-4.86	119.68	126.37
1	1	1917	PSU	C4-N3-C2	-4.85	119.69	126.37
2	2	1207	2MG	C5-C4-N3	-4.79	120.76	128.39
1	1	955	PSU	C4-N3-C2	-4.78	119.79	126.37
1	1	2069	G7M	C5-C4-N3	-4.78	119.12	128.15
2	2	1516	2MG	C2-N1-C6	-4.78	118.78	124.55
2	2	1518	MA6	N9-C8-N7	-4.75	107.20	113.94
1	1	2504	PSU	N1-C2-N3	4.73	120.16	115.17
2	2	527	7MG	C5-C6-N1	4.71	119.24	110.94
2	2	1207	2MG	C2-N1-C6	-4.67	118.90	124.55
52	5	54	5MU	C5M-C5-C4	4.67	123.77	118.78
1	1	2504	PSU	C4-N3-C2	-4.60	120.04	126.37
1	1	747	5MU	N3-C2-N1	4.51	120.77	114.89
1	1	955	PSU	N1-C2-N3	4.48	119.89	115.17
2	2	516	PSU	N1-C2-N3	4.44	119.85	115.17
1	1	2069	G7M	C5-C6-N1	4.43	121.01	111.84
1	1	2503	2MA	N9-C8-N7	-4.41	107.68	113.94
1	1	1835	2MG	N9-C4-N3	4.41	134.76	125.95
1	1	2030	6MZ	C1'-N9-C8	-4.36	117.42	127.09
1	1	2251	OMG	C2-N3-C4	4.34	119.78	112.30
42	q	89	0TD	OD2-CG-CB	4.34	122.52	113.15
1	1	1618	6MZ	C1'-N9-C8	-4.30	117.55	127.09
1	1	2030	6MZ	C5-N7-C8	4.28	110.17	103.45
2	2	1519	MA6	N9-C8-N7	-4.25	107.91	113.94
2	2	527	7MG	C2-N3-C4	4.24	119.61	112.30
2	2	1519	MA6	C5-C4-N3	-4.22	120.91	126.72
52	5	55	PSU	N1-C2-N3	4.15	119.55	115.17
2	2	1518	MA6	C5-C4-N3	-4.15	121.00	126.72
52	5	55	PSU	C4-N3-C2	-4.13	120.69	126.37
1	1	2445	2MG	N9-C4-N3	4.11	134.16	125.95
1	1	2503	2MA	N3-C4-N9	4.04	132.12	126.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	2552	OMU	N3-C2-N1	4.02	120.13	114.89
1	1	2552	OMU	C5-C4-N3	3.95	120.33	114.80
1	1	747	5MU	C5M-C5-C6	-3.90	117.58	122.85
2	2	1519	MA6	C2-N1-C6	3.88	121.31	111.83
1	1	1939	5MU	C5M-C5-C6	-3.85	117.64	122.85
2	2	527	7MG	C5-C4-N3	-3.83	120.93	128.13
1	1	2069	G7M	C2-N1-C6	-3.82	118.17	125.11
1	1	745	1MG	N9-C8-N7	-3.82	106.31	113.40
52	5	54	5MU	N3-C2-N1	3.80	119.84	114.89
2	2	1207	2MG	N2-C2-N3	-3.80	115.67	120.51
52	5	54	5MU	C4-N3-C2	-3.80	122.36	127.34
1	1	2503	2MA	C4-N9-C8	3.78	109.71	105.74
1	1	745	1MG	C2-N3-C4	3.75	120.42	111.98
52	5	54	5MU	C5M-C5-C6	-3.75	117.78	122.85
2	2	1498	UR3	C5-C4-N3	3.74	119.97	115.04
1	1	1915	3TD	C4-N3-C2	-3.72	120.67	124.61
2	2	1518	MA6	C2-N1-C6	3.70	120.86	111.83
1	1	1618	6MZ	C5-N7-C8	3.69	109.25	103.45
2	2	527	7MG	C5-C4-N9	3.61	110.95	106.33
1	1	1939	5MU	O2-C2-N1	-3.59	118.13	122.80
1	1	2457	PSU	O2-C2-N1	-3.59	119.09	122.79
2	2	1207	2MG	CM2-N2-C2	-3.57	115.97	123.65
1	1	2504	PSU	C6-N1-C2	-3.56	119.38	122.69
1	1	2251	OMG	C2-N1-C6	-3.52	118.73	125.11
1	1	1618	6MZ	C5-C4-N3	-3.50	121.90	126.72
1	1	1962	5MC	C5-C6-N1	-3.47	119.54	123.31
1	1	2251	OMG	N9-C4-N3	3.47	132.89	125.95
52	5	8	4SU	N3-C2-N1	3.47	119.41	114.89
1	1	2069	G7M	C1'-N9-C4	-3.46	116.28	126.49
1	1	2457	PSU	C6-C5-C4	3.44	120.49	118.17
1	1	1618	6MZ	C2-N3-C4	3.43	120.21	111.83
1	1	2504	PSU	O2-C2-N1	-3.42	119.26	122.79
1	1	2069	G7M	O6-C6-C5	-3.42	120.38	128.01
2	2	1207	2MG	C1'-N9-C4	-3.41	116.43	126.49
1	1	1911	PSU	O2-C2-N1	-3.39	119.29	122.79
2	2	1519	MA6	C4-C5-C6	3.39	119.41	115.91
1	1	2069	G7M	N9-C4-N3	3.35	132.66	125.95
1	1	2445	2MG	C5-C6-N1	3.34	121.76	113.25
2	2	1516	2MG	N9-C4-N3	3.33	132.61	125.95
1	1	1917	PSU	C6-N1-C2	-3.32	119.61	122.69
1	1	2030	6MZ	C2-N3-C4	3.31	119.91	111.83
2	2	1519	MA6	C2-N3-C4	3.28	119.85	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	5	20	H2U	C5-C6-N1	3.27	121.42	111.52
2	2	1518	MA6	C2-N3-C4	3.26	119.79	111.83
2	2	1518	MA6	C4-C5-C6	3.25	119.28	115.91
1	1	2251	OMG	N9-C8-N7	-3.25	107.37	113.40
2	2	527	7MG	C4-C5-N7	3.21	109.17	105.38
52	5	8	4SU	C5-C4-S4	-3.19	120.66	124.31
2	2	1207	2MG	C1'-N9-C8	3.17	135.74	126.73
1	1	747	5MU	C6-C5-C4	3.17	120.63	118.02
2	2	967	5MC	C5-C6-N1	-3.15	119.89	123.31
2	2	1518	MA6	C4-N9-C8	3.12	109.02	105.74
1	1	1962	5MC	C5-C4-N4	-3.12	117.00	121.39
1	1	2030	6MZ	C5-C4-N3	-3.12	122.42	126.72
2	2	1407	5MC	O2-C2-N3	-3.11	117.43	122.33
1	1	1939	5MU	C5M-C5-C4	3.10	122.10	118.78
1	1	2580	PSU	O2-C2-N1	-3.10	119.59	122.79
52	5	20	H2U	N3-C2-N1	3.10	119.76	116.65
2	2	1518	MA6	C5-N7-C8	3.07	108.27	103.45
1	1	746	PSU	O2-C2-N1	-3.06	119.64	122.79
2	2	1519	MA6	C5-N7-C8	3.05	108.25	103.45
1	1	2457	PSU	C6-N1-C2	-3.01	119.90	122.69
1	1	1917	PSU	O2-C2-N1	-3.00	119.69	122.79
52	5	8	4SU	C1'-N1-C2	2.98	122.95	117.59
1	1	1911	PSU	C6-N1-C2	-2.97	119.94	122.69
2	2	1516	2MG	C1'-N9-C4	-2.97	117.73	126.49
1	1	2503	2MA	N6-C6-N1	2.96	121.02	117.03
1	1	1835	2MG	C5-C6-N1	2.95	120.77	113.25
2	2	1407	5MC	C5-C6-N1	-2.94	120.12	123.31
1	1	2552	OMU	O4-C4-C5	-2.92	120.12	125.16
2	2	966	2MG	C5-C6-N1	2.89	120.60	113.25
1	1	745	1MG	C5-C6-N1	2.88	120.34	115.02
1	1	2445	2MG	N9-C8-N7	-2.86	108.10	113.40
1	1	745	1MG	N9-C4-N3	2.85	131.66	125.95
1	1	2445	2MG	O6-C6-C5	-2.85	119.01	126.53
2	2	1516	2MG	O6-C6-C5	-2.85	119.02	126.53
2	2	1498	UR3	C1'-N1-C2	2.84	121.69	117.04
1	1	2552	OMU	O2-C2-N1	-2.84	119.11	122.80
1	1	2030	6MZ	C4-C5-C6	-2.83	114.42	116.78
1	1	747	5MU	C5M-C5-C4	2.82	121.79	118.78
2	2	1207	2MG	N9-C8-N7	-2.78	108.25	113.40
1	1	745	1MG	N2-C2-N1	2.76	121.01	118.79
2	2	1516	2MG	C5-C6-N1	2.75	120.25	113.25
2	2	1402	4OC	O2-C2-N3	-2.74	118.01	122.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	1939	5MU	O4-C4-N3	-2.74	114.97	120.11
2	2	966	2MG	O6-C6-C5	-2.73	119.34	126.53
1	1	2251	OMG	C5-C6-N1	2.72	120.19	113.25
1	1	2445	2MG	CM2-N2-C2	-2.72	117.80	123.65
1	1	1939	5MU	C6-C5-C4	2.72	120.26	118.02
1	1	1835	2MG	N9-C8-N7	-2.72	108.36	113.40
2	2	1516	2MG	N9-C8-N7	-2.72	108.36	113.40
42	q	89	0TD	OD1-CG-CB	-2.71	116.77	122.44
1	1	2069	G7M	C1'-N9-C8	2.70	135.87	126.74
1	1	1962	5MC	C1'-N1-C6	-2.69	116.72	121.15
1	1	2580	PSU	C6-N1-C2	-2.67	120.21	122.69
1	1	1618	6MZ	C4-C5-C6	-2.64	114.58	116.78
52	5	20	H2U	C5-C4-N3	2.64	119.50	116.69
1	1	2030	6MZ	C6-C5-N7	2.64	135.31	132.43
2	2	1519	MA6	C4-C5-N7	-2.64	107.56	110.58
2	2	1516	2MG	C1'-N9-C8	2.59	134.10	126.73
1	1	2251	OMG	O6-C6-C5	-2.57	119.74	126.53
2	2	1207	2MG	O6-C6-C5	-2.56	119.78	126.53
52	5	54	5MU	C1'-N1-C2	2.56	122.18	117.59
52	5	55	PSU	C6-N1-C2	-2.55	120.33	122.69
2	2	1518	MA6	C4-C5-N7	-2.51	107.71	110.58
2	2	527	7MG	C2-N1-C6	-2.50	120.57	125.11
2	2	1207	2MG	C5-C6-N1	2.50	119.61	113.25
52	5	54	5MU	O4-C4-N3	-2.44	115.52	120.11
1	1	1835	2MG	O6-C6-C5	-2.44	120.08	126.53
1	1	955	PSU	O2-C2-N1	-2.43	120.28	122.79
2	2	1207	2MG	N9-C4-N3	2.43	130.81	125.95
1	1	745	1MG	C8-N7-C5	2.41	108.55	104.26
1	1	2605	PSU	C6-C5-C4	2.41	119.80	118.17
1	1	747	5MU	C1'-N1-C6	-2.38	117.24	121.15
2	2	1518	MA6	N3-C4-N9	2.38	131.21	127.17
2	2	966	2MG	N9-C8-N7	-2.37	109.00	113.40
1	1	2503	2MA	N3-C2-N1	-2.34	121.63	125.77
1	1	955	PSU	C6-N1-C2	-2.34	120.52	122.69
1	1	1835	2MG	CM2-N2-C2	-2.31	118.69	123.65
2	2	527	7MG	N9-C8-N7	2.30	106.63	103.37
52	5	32	4OC	O2-C2-N3	-2.29	118.72	122.33
52	5	32	4OC	C6-C5-C4	2.29	119.76	117.00
2	2	1519	MA6	C5-C4-N9	2.28	108.29	105.81
2	2	1407	5MC	C5-C4-N3	-2.27	119.43	121.75
2	2	1516	2MG	N1-C2-N2	2.26	118.87	116.56
1	1	746	PSU	C6-C5-C4	2.24	119.69	118.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	516	PSU	O3'-C3'-C4'	2.23	117.50	111.08
1	1	746	PSU	O4-C4-C5	-2.23	118.46	124.01
1	1	2445	2MG	C1'-N9-C4	-2.23	119.89	126.49
1	1	747	5MU	C1'-N1-C2	2.23	121.60	117.59
2	2	1516	2MG	CM2-N2-C2	-2.22	118.87	123.65
1	1	2580	PSU	O4'-C1'-C2'	2.20	108.20	105.15
52	5	55	PSU	O2-C2-N1	-2.20	120.52	122.79
1	1	2503	2MA	C6-C5-C4	2.20	120.18	117.18
1	1	747	5MU	O2-C2-N1	-2.19	119.94	122.80
1	1	2605	PSU	C6-N1-C2	-2.17	120.67	122.69
2	2	967	5MC	O2-C2-N3	-2.17	118.91	122.33
1	1	2605	PSU	O2-C2-N1	-2.17	120.55	122.79
1	1	2030	6MZ	C4-N9-C1'	-2.17	121.56	126.63
1	1	2503	2MA	C5-N7-C8	2.16	106.84	103.45
2	2	516	PSU	C6-N1-C2	-2.16	120.69	122.69
2	2	1519	MA6	C4-N9-C8	2.14	107.98	105.74
2	2	1519	MA6	N3-C4-N9	2.13	130.80	127.17
2	2	967	5MC	CM5-C5-C6	-2.13	119.96	122.85
1	1	2251	OMG	C8-N7-C5	2.09	107.99	104.26
53	H	77	FME	O1-CN-N	-2.09	119.91	125.32
1	1	2457	PSU	O4'-C1'-C2'	2.09	108.04	105.15
2	2	1402	4OC	CM4-N4-C4	-2.08	118.39	122.45
2	2	1402	4OC	C6-C5-C4	2.06	119.48	117.00
1	1	2503	2MA	CM2-C2-N1	2.05	120.20	117.13
1	1	1835	2MG	N1-C2-N2	2.05	118.65	116.56
1	1	2605	PSU	C5-C6-N1	-2.04	119.30	122.14
1	1	747	5MU	O4-C4-N3	-2.03	116.29	120.11
1	1	746	PSU	C6-N1-C2	-2.02	120.82	122.69
2	2	527	7MG	O6-C6-C5	-2.01	122.69	127.62

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	1	746	PSU	C2'-C1'-C5-C4
1	1	746	PSU	C2'-C1'-C5-C6
1	1	1618	6MZ	C5-C6-N6-C9
1	1	1618	6MZ	N1-C6-N6-C9
1	1	1915	3TD	O4'-C1'-C5-C4
1	1	1915	3TD	O4'-C1'-C5-C6
1	1	2030	6MZ	C3'-C4'-C5'-O5'
2	2	516	PSU	O4'-C1'-C5-C4

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Mol	Chain	Res	Type	Atoms
2	2	516	PSU	O4'-C1'-C5-C6
2	2	1519	MA6	O4'-C4'-C5'-O5'
52	5	20	H2U	O4'-C1'-N1-C2
52	5	20	H2U	O4'-C1'-N1-C6
53	H	77	FME	O1-CN-N-CA
1	1	2445	2MG	C3'-C4'-C5'-O5'
2	2	527	7MG	C3'-C4'-C5'-O5'
2	2	1402	4OC	O4'-C4'-C5'-O5'
2	2	1519	MA6	C3'-C4'-C5'-O5'
52	5	8	4SU	O4'-C1'-N1-C2
1	1	2030	6MZ	O4'-C4'-C5'-O5'
1	1	2445	2MG	O4'-C4'-C5'-O5'
2	2	527	7MG	O4'-C4'-C5'-O5'
52	5	8	4SU	C3'-C4'-C5'-O5'
52	5	20	H2U	O4'-C4'-C5'-O5'
52	5	20	H2U	C3'-C4'-C5'-O5'
1	1	2503	2MA	O4'-C4'-C5'-O5'
52	5	55	PSU	C3'-C4'-C5'-O5'
53	H	77	FME	CA-CB-CG-SD
2	2	966	2MG	O4'-C4'-C5'-O5'
2	2	966	2MG	C3'-C4'-C5'-O5'
2	2	1402	4OC	C3'-C4'-C5'-O5'
52	5	8	4SU	O4'-C1'-N1-C6
52	5	55	PSU	O4'-C4'-C5'-O5'
42	q	89	0TD	CG-CB-SB-CSB
1	1	1962	5MC	C2'-C1'-N1-C6
52	5	8	4SU	O4'-C4'-C5'-O5'
2	2	1519	MA6	C4'-C5'-O5'-P
1	1	746	PSU	O4'-C1'-C5-C4
1	1	2457	PSU	O4'-C1'-C5-C4
1	1	747	5MU	C4'-C5'-O5'-P
2	2	527	7MG	C4'-C5'-O5'-P
2	2	966	2MG	C4'-C5'-O5'-P
1	1	1915	3TD	O4'-C4'-C5'-O5'
1	1	1962	5MC	O4'-C1'-N1-C6
1	1	746	PSU	O4'-C1'-C5-C6
1	1	2457	PSU	O4'-C1'-C5-C6
1	1	2503	2MA	O4'-C1'-N9-C8
1	1	1962	5MC	C2'-C1'-N1-C2
1	1	2503	2MA	C3'-C4'-C5'-O5'
1	1	2504	PSU	O4'-C4'-C5'-O5'

There are no ring outliers.

12 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	1	2030	6MZ	1	0
2	2	1519	MA6	1	0
42	q	89	0TD	1	0
2	2	527	7MG	1	0
2	2	1407	5MC	1	0
52	5	8	4SU	1	0
1	1	747	5MU	1	0
1	1	1962	5MC	1	0
2	2	1516	2MG	1	0
52	5	55	PSU	1	0
1	1	1939	5MU	1	0
1	1	2503	2MA	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 296 ligands modelled in this entry, 296 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

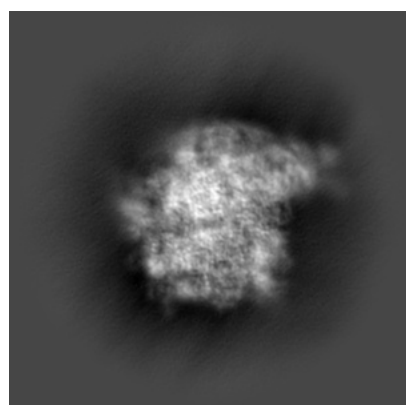
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20174. These allow visual inspection of the internal detail of the map and identification of artifacts.

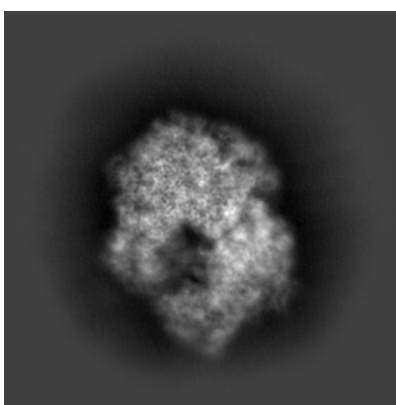
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

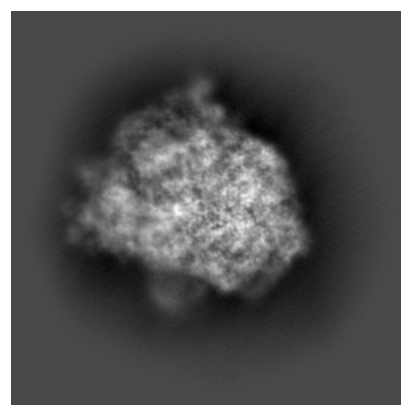
6.1.1 Primary map



X



Y

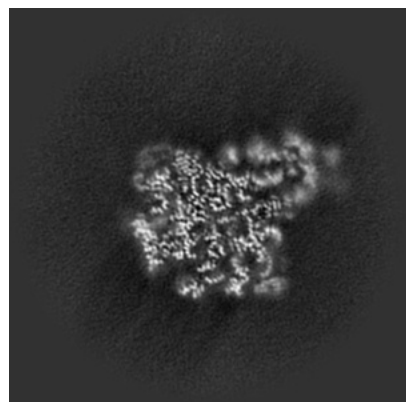


Z

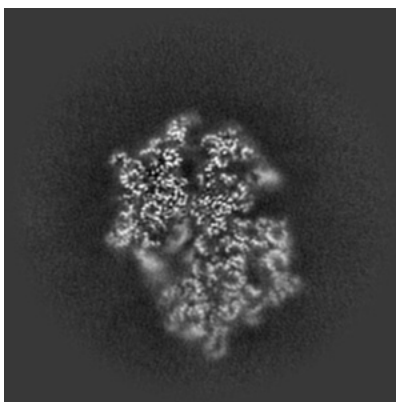
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

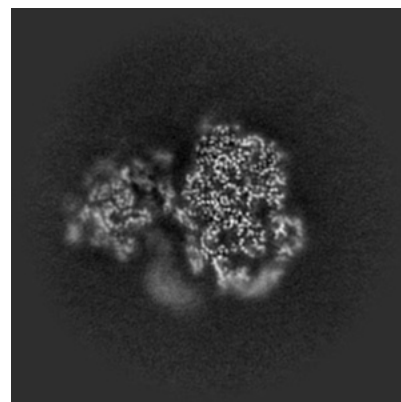
6.2.1 Primary map



X Index: 128



Y Index: 128

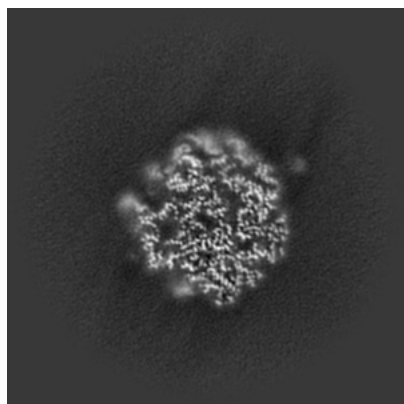


Z Index: 128

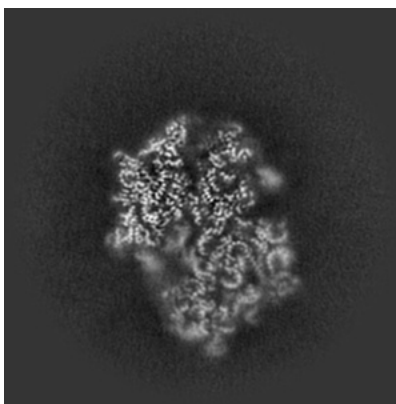
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

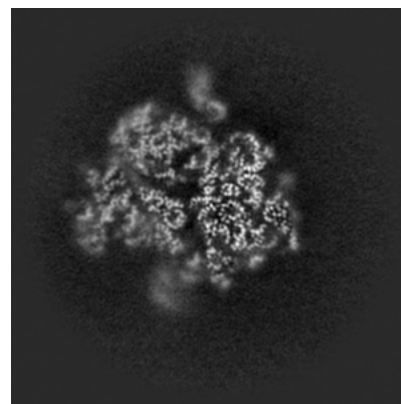
6.3.1 Primary map



X Index: 148



Y Index: 126

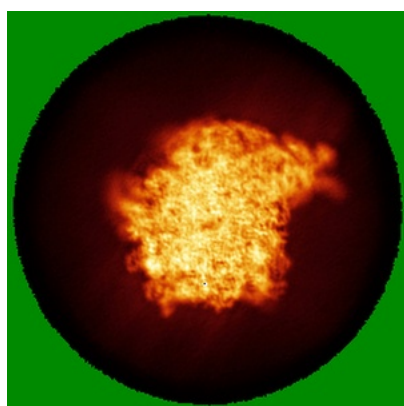


Z Index: 142

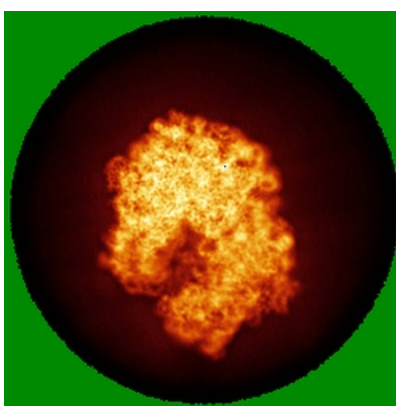
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

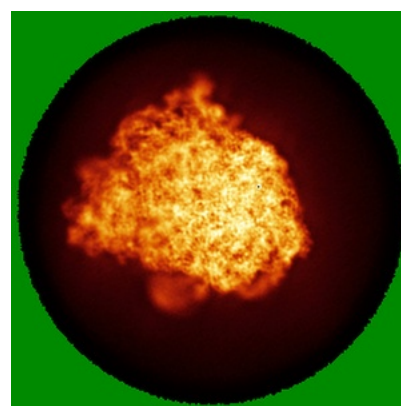
6.4.1 Primary map



X



Y

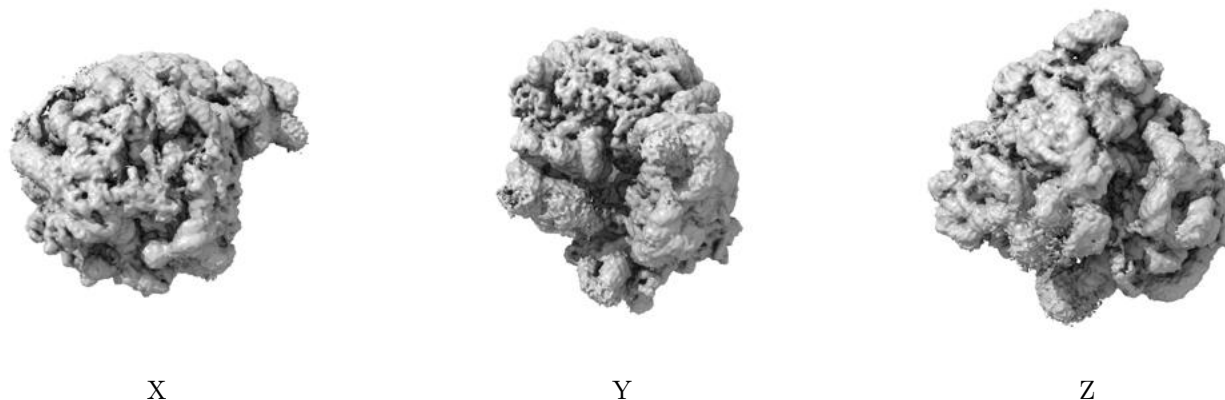


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

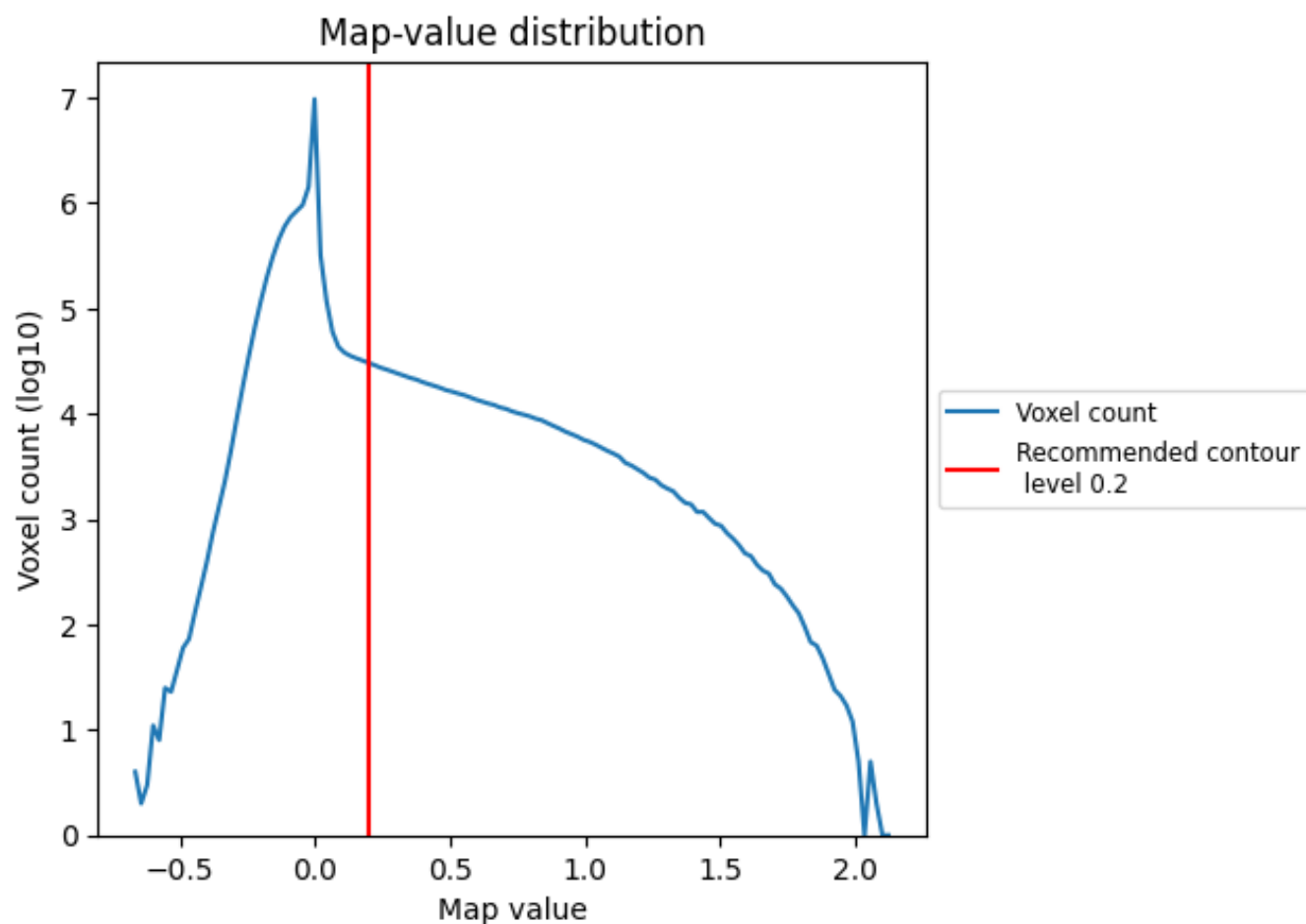
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

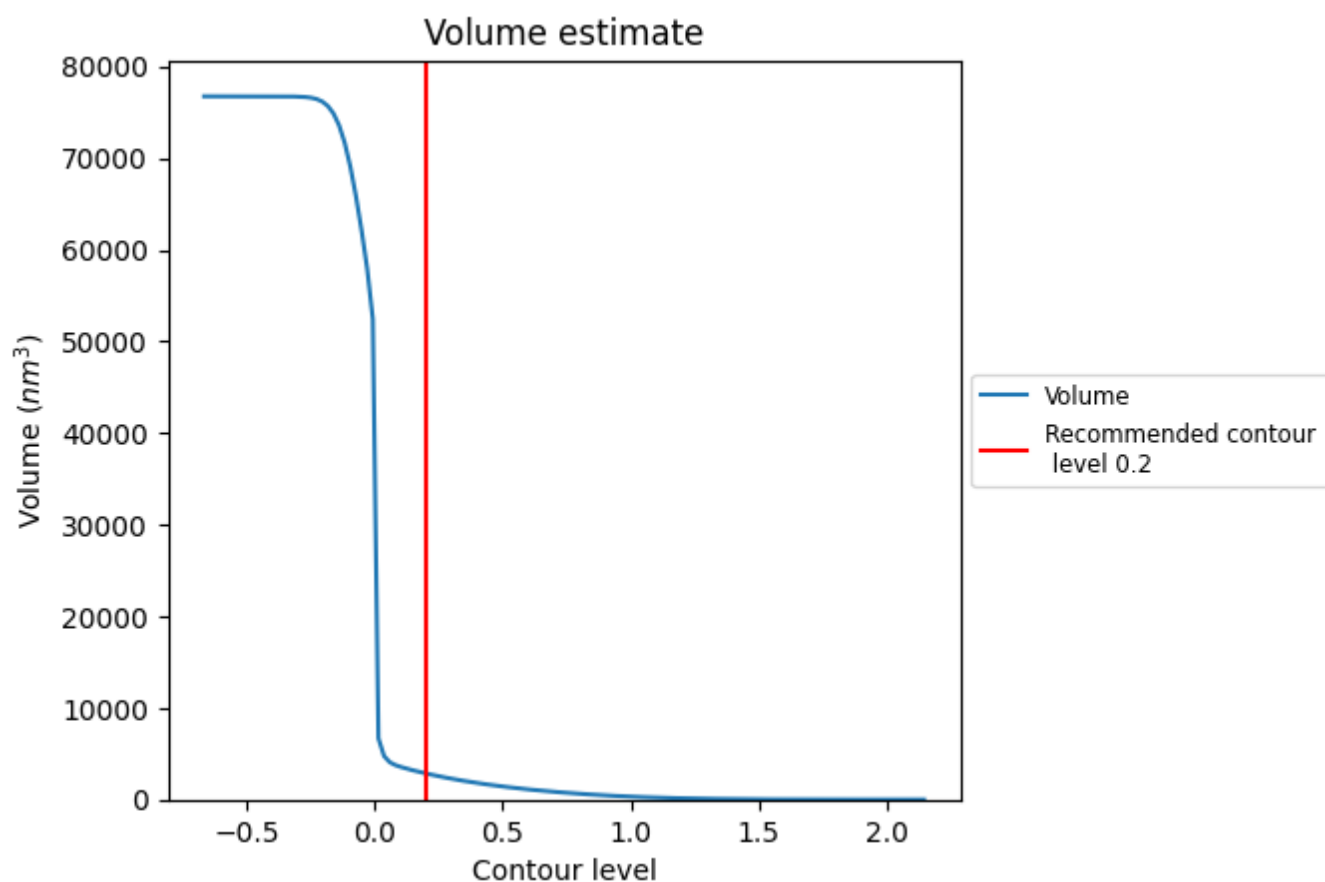
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

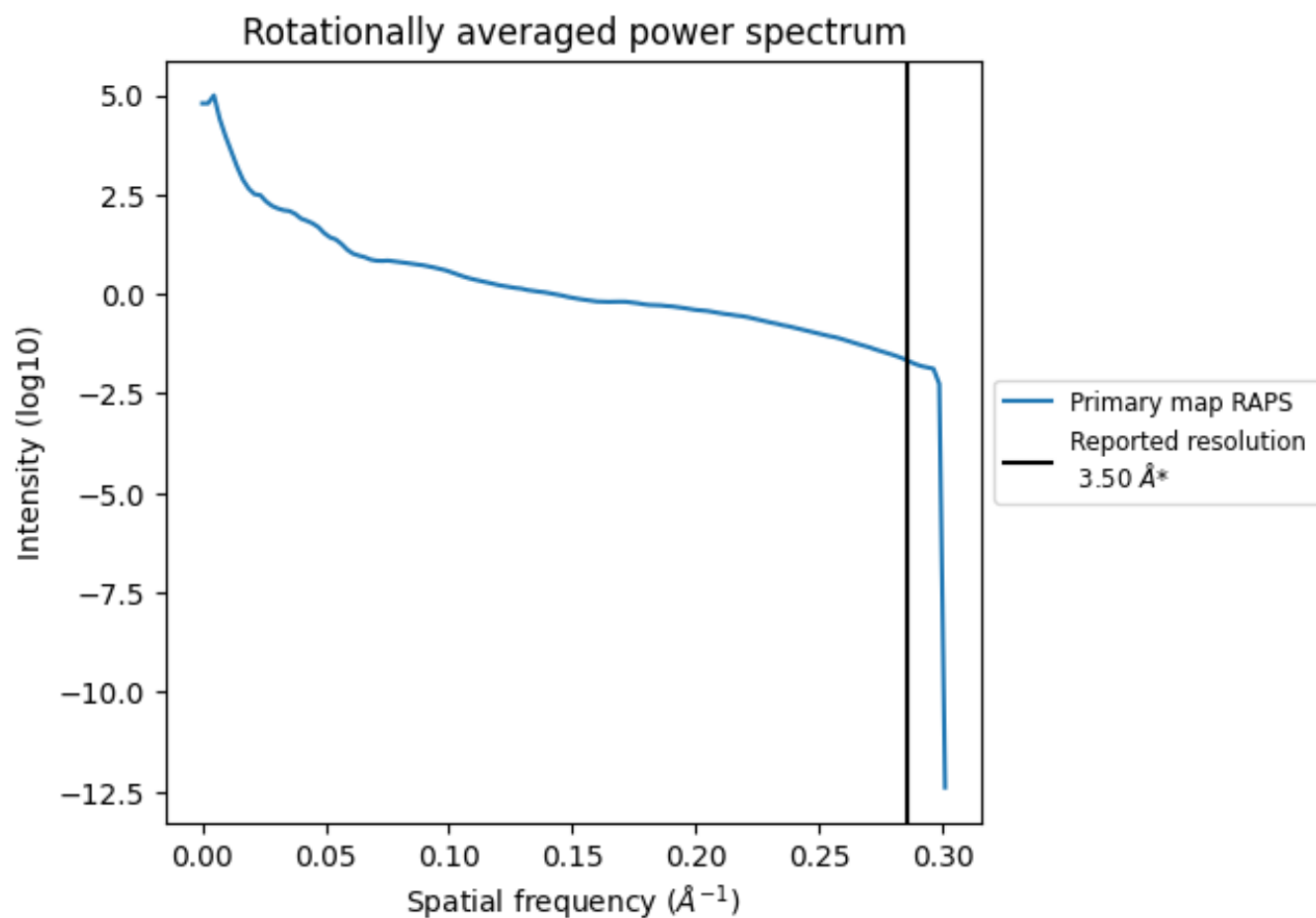
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2849 nm³; this corresponds to an approximate mass of 2574 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

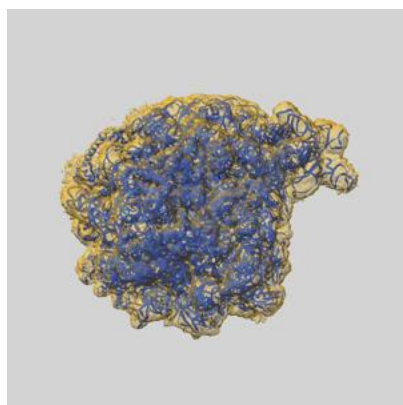
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

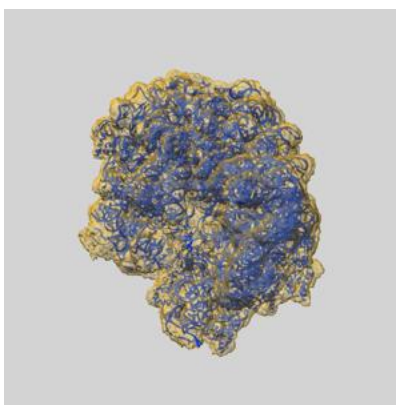
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20174 and PDB model 6ORL. Per-residue inclusion information can be found in section 3 on page 15.

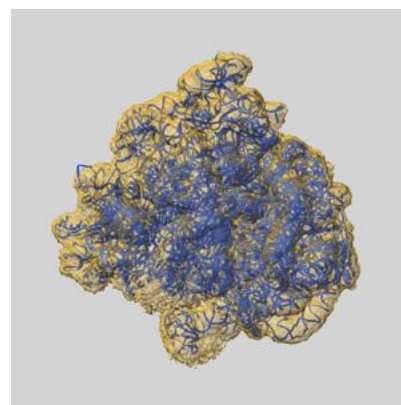
9.1 Map-model overlay [i](#)



X



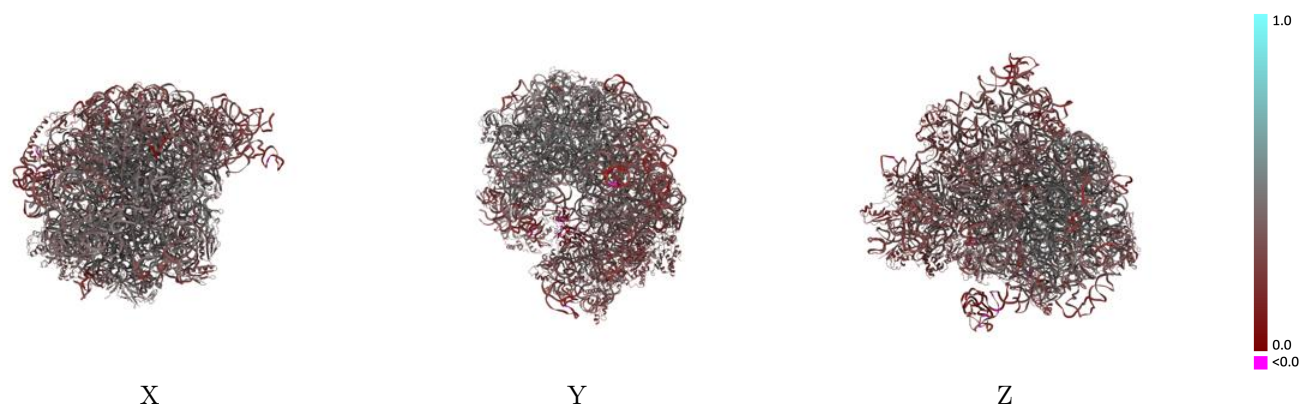
Y



Z

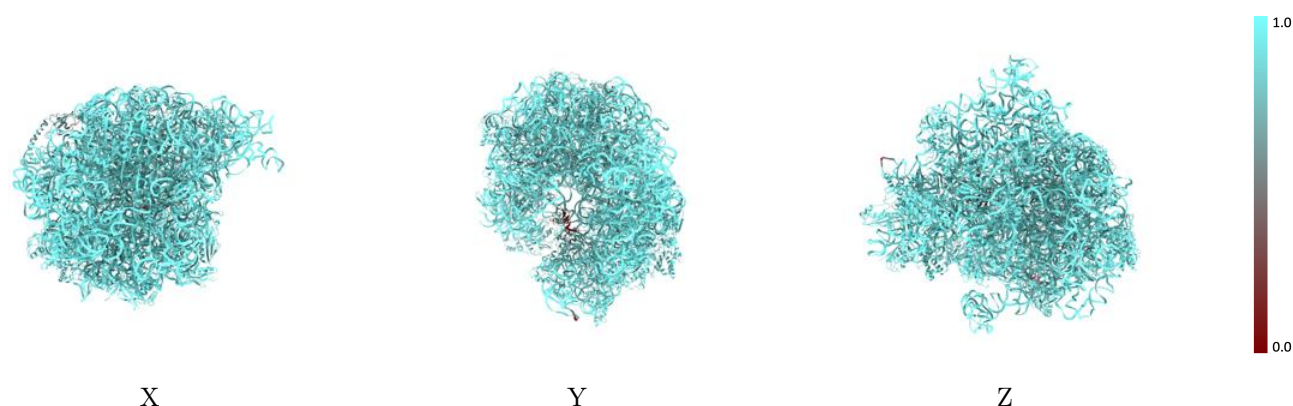
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



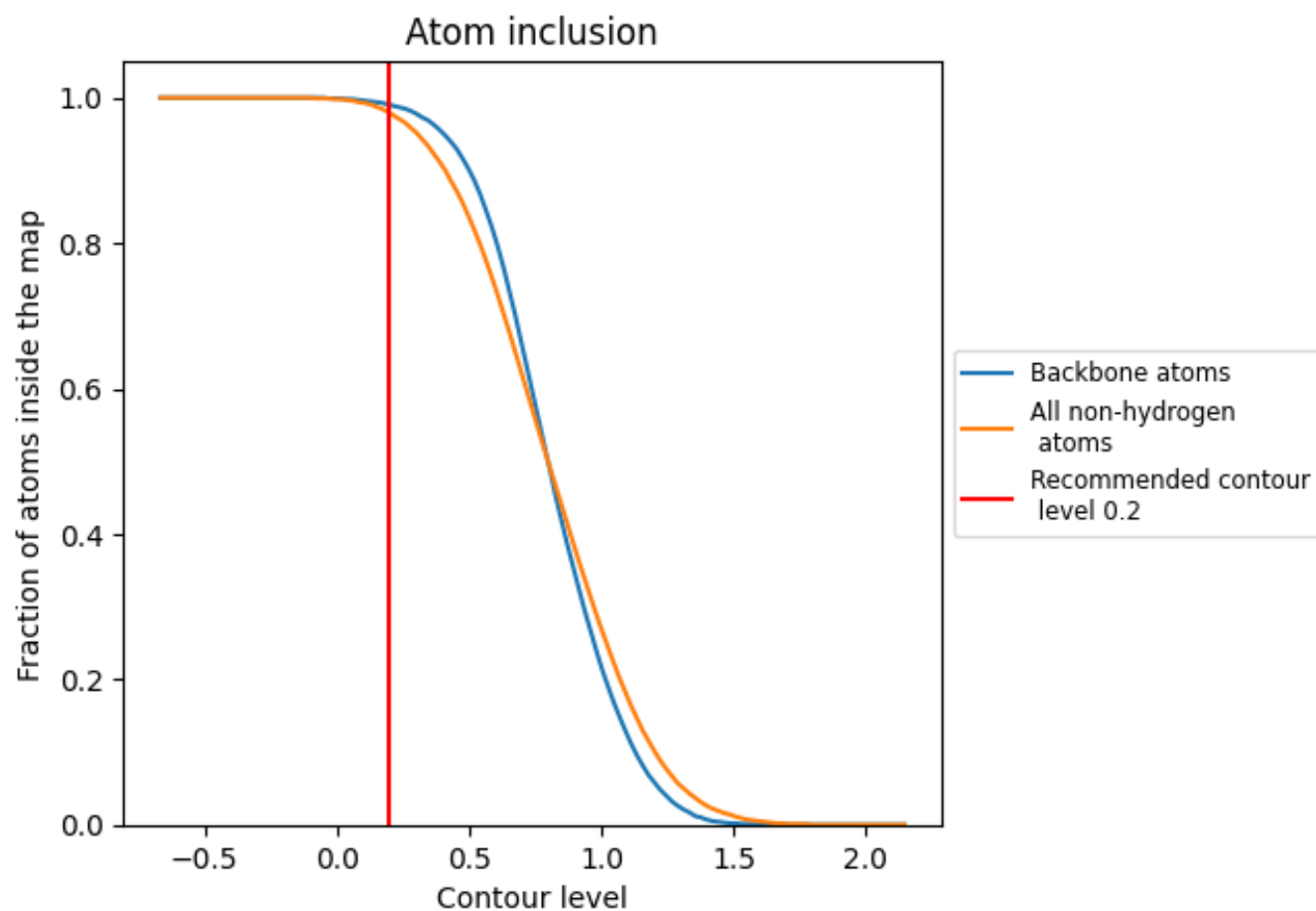
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 98% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9790	 0.3700
1	 0.9990	 0.3980
2	 0.9970	 0.3440
3	 0.9990	 0.3330
4	 1.0000	 0.3850
5	 0.9630	 0.2890
A	 0.5300	 0.1470
B	 0.9620	 0.4470
C	 0.9680	 0.4550
D	 0.9490	 0.4140
E	 0.9370	 0.2860
F	 0.9810	 0.3490
G	 0.7100	 0.2710
H	 0.9380	 0.2890
J	 0.9620	 0.4370
K	 0.9160	 0.4320
L	 0.9770	 0.4340
M	 0.9540	 0.4240
N	 0.9810	 0.4340
O	 0.9510	 0.3320
P	 0.9510	 0.4300
Q	 0.9760	 0.4280
R	 0.9550	 0.4400
S	 0.9590	 0.4450
T	 0.9470	 0.4050
U	 0.9640	 0.3850
V	 0.9770	 0.3640
W	 0.9540	 0.4400
X	 0.9620	 0.4050
Y	 0.9610	 0.3160
Z	 0.9310	 0.4230
b	 0.9770	 0.4290
c	 0.9710	 0.3990
d	 0.9610	 0.4380
e	 0.9330	 0.4330



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Chain	Atom inclusion	Q-score
f	 0.9630	 0.4030
g	 0.9560	 0.3030
h	 0.9540	 0.3380
i	 0.9740	 0.2700
j	 0.9770	 0.3610
k	 0.9720	 0.3300
l	 0.9560	 0.2770
m	 0.9740	 0.3540
n	 0.9850	 0.2890
o	 0.9580	 0.2920
p	 0.9840	 0.3530
q	 0.9170	 0.3580
r	 0.9450	 0.2770
s	 0.9760	 0.3240
t	 0.9680	 0.3250
u	 0.9750	 0.3400
v	 0.9680	 0.3120
w	 0.9920	 0.3140
x	 0.9720	 0.2870
y	 0.9450	 0.2680
z	 0.9610	 0.2880