



Full wwPDB EM Validation Report ⓘ

Jun 23, 2026 – 11:50 AM JST

PDB ID : 8HTZ / pdb_00008htz
EMDB ID : EMD-35020
Title : E. coli 70S ribosome complexed with H. marismortui tRNA_Ile2 bearing agm2C34 in classical state
Authors : Akiyama, N.; Ishiguro, K.; Yokoyama, T.; Shirouzu, M.; Suzuki, T.
Deposited on : 2022-12-22
Resolution : 2.40 Å(reported)
Based on initial model : 7K00

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 : 1.8.5 (274361), CSD as541be (2020)
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

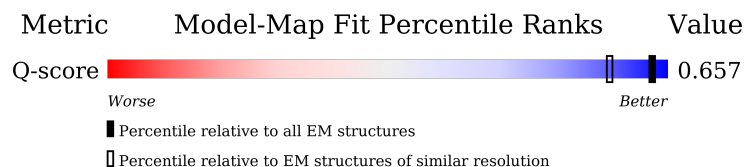
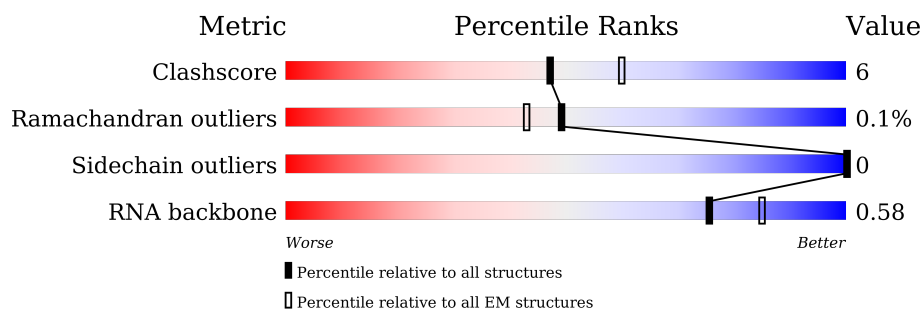
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.40 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	5628 (1.90 - 2.90)

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	A	1542	
2	B	241	
3	C	233	

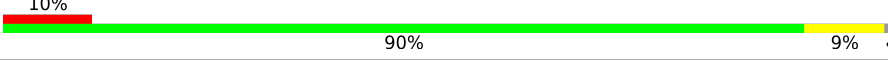
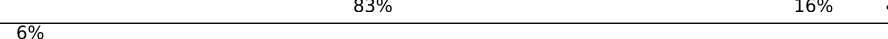
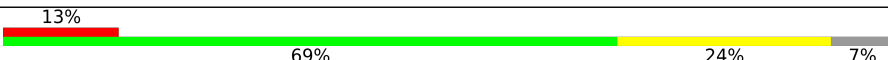
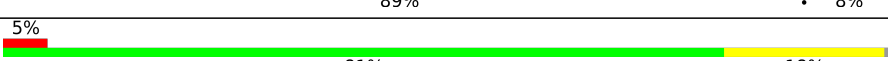
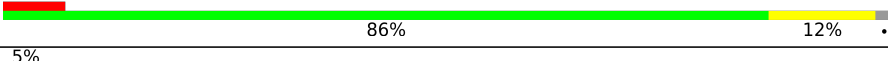
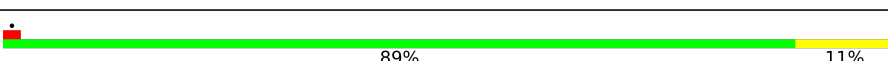
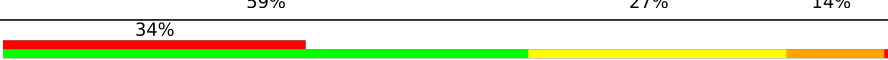
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Mol	Chain	Length	Quality of chain
4	D	206	
5	E	167	
6	F	135	
7	G	179	
8	H	130	
9	I	130	
10	J	103	
11	K	129	
12	L	124	
13	M	118	
14	N	101	
15	O	89	
16	P	82	
17	Q	84	
18	R	75	
19	S	92	
20	T	87	
21	U	71	
22	a	2904	
23	b	120	
24	c	273	
25	d	209	
26	e	201	
27	f	179	
28	g	177	

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Mol	Chain	Length	Quality of chain
29	h	149	
30	i	142	
31	j	123	
32	k	144	
33	l	136	
34	m	127	
35	n	117	
36	o	115	
37	p	118	
38	q	103	
39	r	110	
40	s	100	
41	t	104	
42	u	94	
43	v	85	
44	w	78	
45	x	63	
46	y	59	
47	z	57	
48	0	55	
49	1	46	
50	2	65	
51	3	38	
52	4	70	
53	V	76	

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Mol	Chain	Length	Quality of chain
54	W	48	
55	X	78	

2 Entry composition

There are 57 unique types of molecules in this entry. The entry contains 141919 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 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1519	Total	C	N	O	P	0	0
			32612	14552	5986	10555	1519		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	224	Total	C	N	O	S	0	0
			1753	1109	315	321	8		

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

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

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

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

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

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

- Molecule 6 is a protein called 30S ribosomal protein S6, fully modified isoform.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	103	Total	C	N	O	S	0	0
			839	530	151	151	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	153	Total	C	N	O	S	0	0
			1203	750	231	218	4		

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	98	Total	C	N	O	S	0	0
			786	493	150	142	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	117	Total	C	N	O	S	0	0
			877	540	173	161	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	119	IAS	ASN	modified residue	UNP P0A7R9

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	115	Total	C	N	O	S	0	0
			891	552	179	157	3		

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	81	Total	C	N	O	S	0	0
			643	403	127	112	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	79	Total	C	N	O	S	0	0
			641	406	120	112	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
18	R	66	Total	C	N	O	S	0	0
			544	345	102	96	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
19	S	84	Total	C	N	O	S	0	0
			668	427	127	112	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	86	Total	C	N	O	S	0	0
			670	414	138	115	3		

- Molecule 21 is a protein called 30S ribosomal protein S21.

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

- Molecule 22 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	a	2753	Total	C	N	O	P	0	0
			59130	26384	10897	19096	2753		

- Molecule 23 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	b	119	Total	C	N	O	P	0	0
			2549	1135	466	829	119		

- Molecule 24 is a protein called 50S ribosomal protein L2.

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

- Molecule 25 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	d	209	Total	C	N	O	S	0	0
			1566	980	288	294	4		

- Molecule 26 is a protein called 50S ribosomal protein L4.

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

- Molecule 27 is a protein called 50S ribosomal protein L5.

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

- Molecule 28 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	g	176	Total	C	N	O	S	0	0
			1323	832	243	246	2		

- Molecule 29 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	h	41	Total	C	N	O	S	0	0
			303	194	54	54	1		

- Molecule 30 is a protein called 50S ribosomal protein L13.

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

- Molecule 31 is a protein called 50S ribosomal protein L14.

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

- Molecule 32 is a protein called 50S ribosomal protein L15.

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

- Molecule 33 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	l	136	Total	C	N	O	S	0	0
			1075	686	205	177	7		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	82	MS6	MET	modified residue	UNP P0ADY7

- Molecule 34 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	m	118	Total	C	N	O	S	0	0
			945	585	194	161	5		

- Molecule 35 is a protein called 50S ribosomal protein L18.

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

- Molecule 36 is a protein called 50S ribosomal protein L19.

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

- Molecule 37 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	p	117	Total	C	N	O		0	0
			947	604	192	151			

- Molecule 38 is a protein called 50S ribosomal protein L21.

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

- Molecule 39 is a protein called 50S ribosomal protein L22.

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

- Molecule 40 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	s	93	Total	C	N	O	S	0	0
			738	466	139	131	2		

- Molecule 41 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	t	102	Total	C	N	O	S	0	0
			779	492	146	141			

- Molecule 42 is a protein called 50S ribosomal protein L25.

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

- Molecule 43 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	v	78	Total	C	N	O	S	0	0
			586	362	116	107	1		

- Molecule 44 is a protein called 50S ribosomal protein L28.

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

- Molecule 45 is a protein called 50S ribosomal protein L29.

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

- Molecule 46 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	y	58	Total	C	N	O	S	0	0
			449	281	87	79	2		

- Molecule 47 is a protein called 50S ribosomal protein L32.

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

- Molecule 48 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	0	51	Total	C	N	O		0	0
			417	269	76	72			

- Molecule 49 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	1	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

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

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

- Molecule 51 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	3	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 52 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	4	60	Total	C	N	O	S	0	0
			480	299	90	85	6		

- Molecule 53 is a RNA chain called P-site tRNA_Glu.

Mol	Chain	Residues	Atoms						AltConf	Trace
53	V	76	Total	C	N	O	P	S	0	0
			1619	724	288	530	76	1		

- Molecule 54 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	W	11	Total	C	N	O	P	0	0
			236	106	44	75	11		

- Molecule 55 is a RNA chain called A-site tRNA_Ile2.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	X	78	Total	C	N	O	P	0	0
			1687	759	300	550	78		

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

Mol	Chain	Residues	Atoms		AltConf
56	A	94	Total	Mg	0
			94	94	
56	N	1	Total	Mg	0
			1	1	
56	a	208	Total	Mg	0
			208	208	
56	b	5	Total	Mg	0
			5	5	
56	d	1	Total	Mg	0
			1	1	
56	p	1	Total	Mg	0
			1	1	
56	z	1	Total	Mg	0
			1	1	
56	V	2	Total	Mg	0
			2	2	
56	W	2	Total	Mg	0
			2	2	
56	X	2	Total	Mg	0
			2	2	

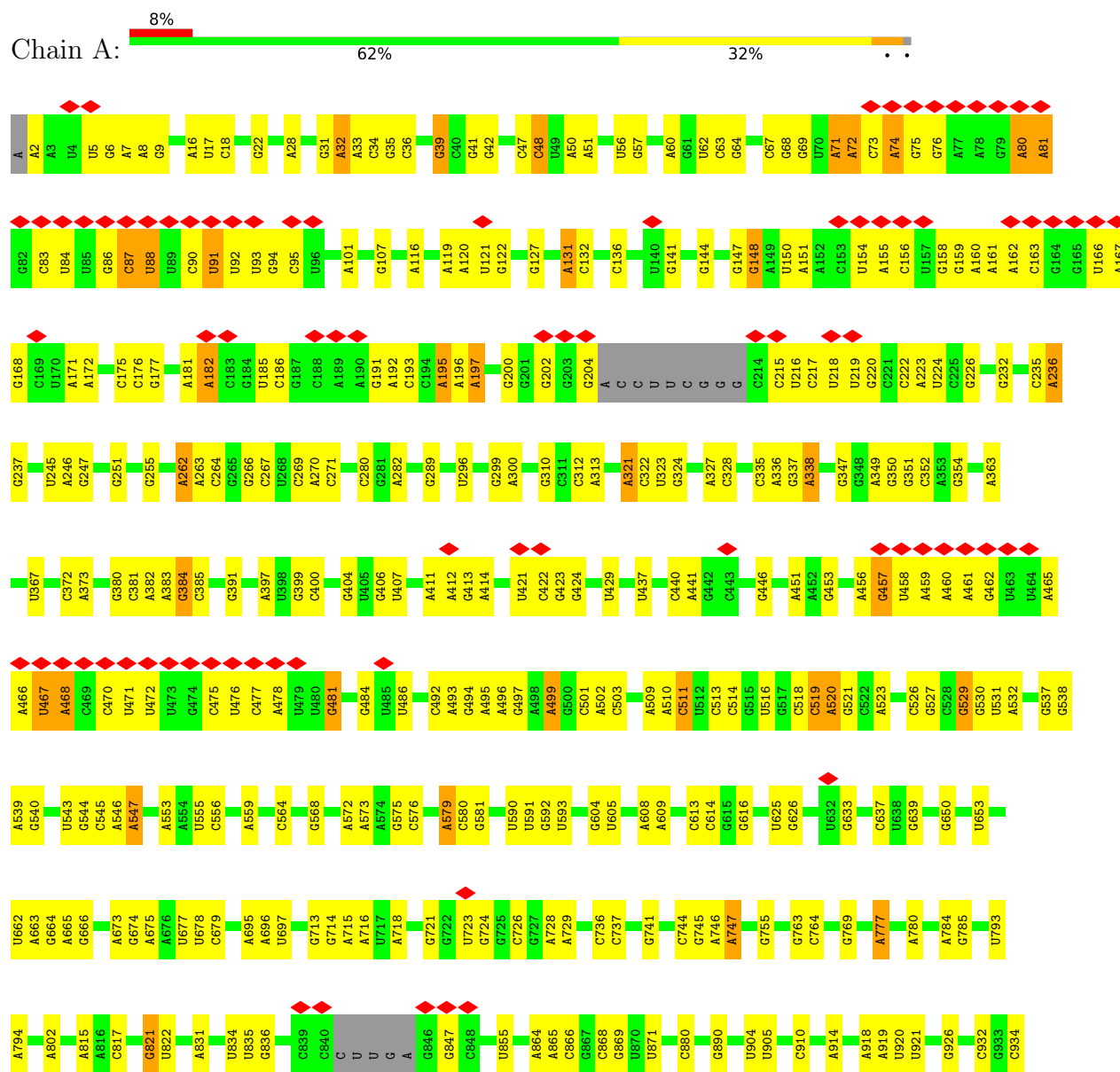
- Molecule 57 is POTASSIUM ION (CCD ID: K) (formula: K).

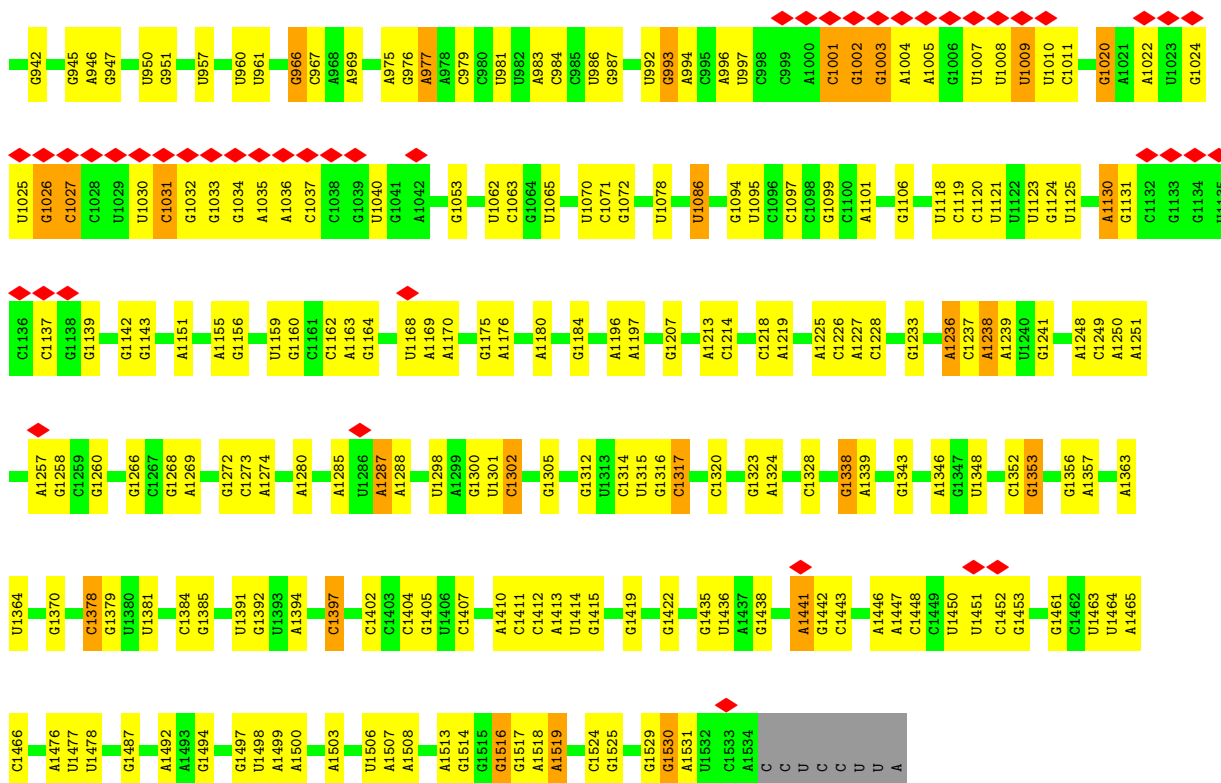
Mol	Chain	Residues	Atoms		AltConf
57	A	1	Total	K	0
			1	1	

3 Residue-property plots

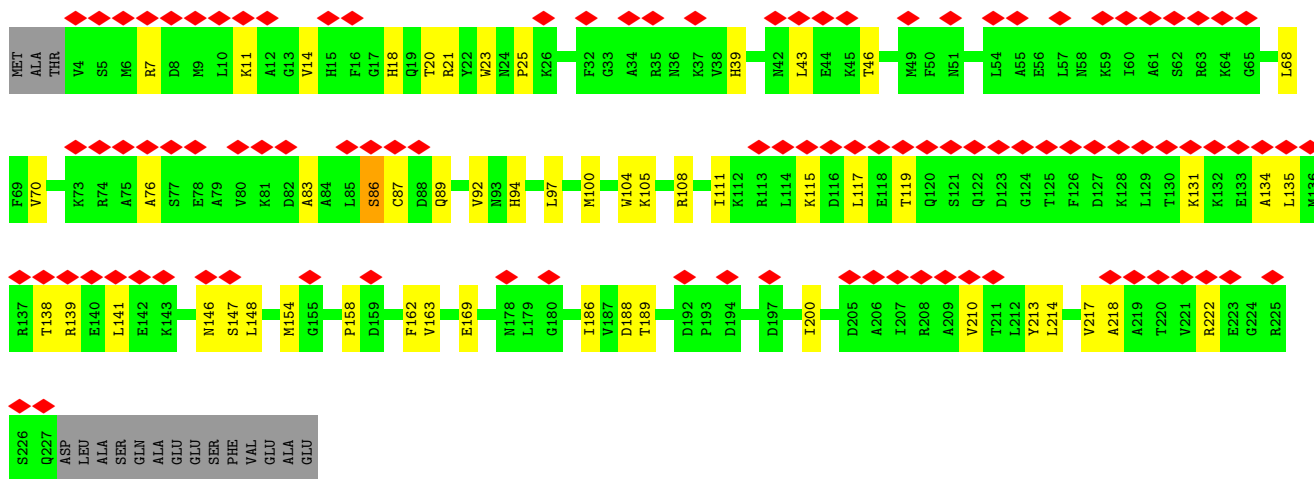
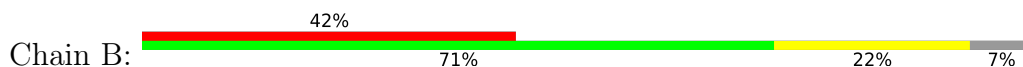
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: 16S rRNA

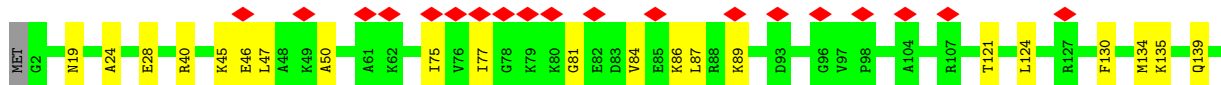
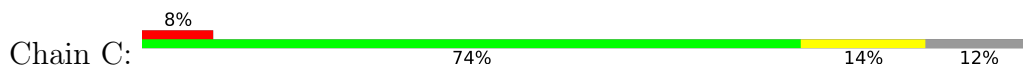


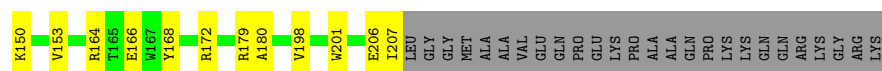


• Molecule 2: 30S ribosomal protein S2

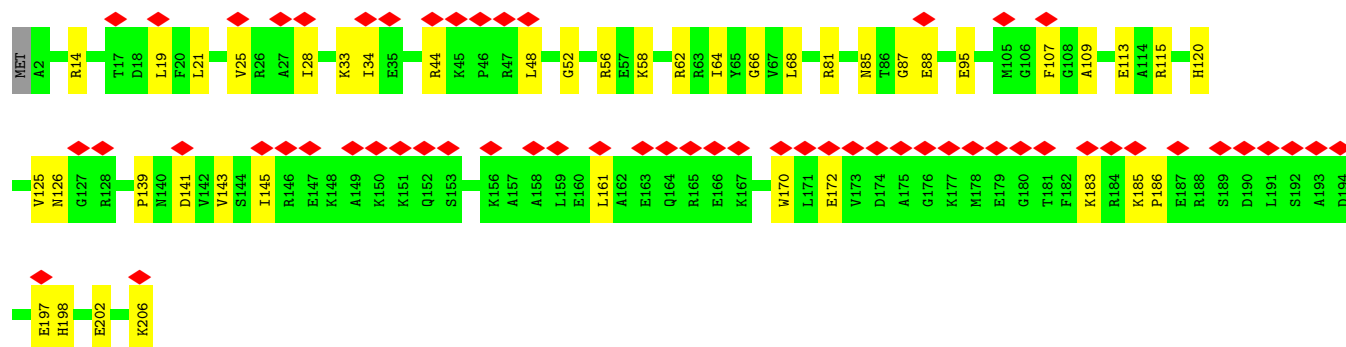
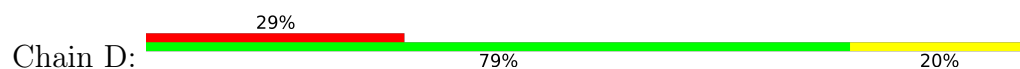


• Molecule 3: 30S ribosomal protein S3

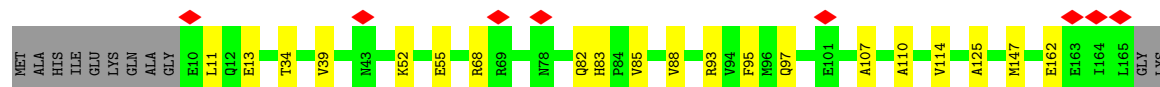
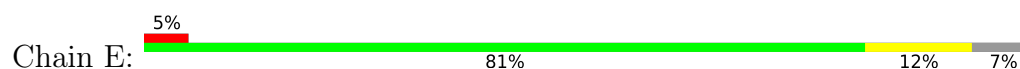




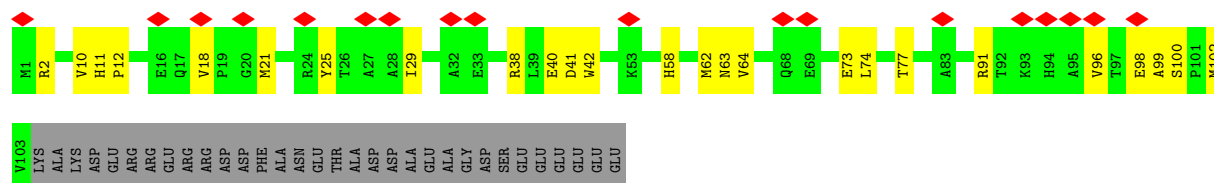
- Molecule 4: 30S ribosomal protein S4



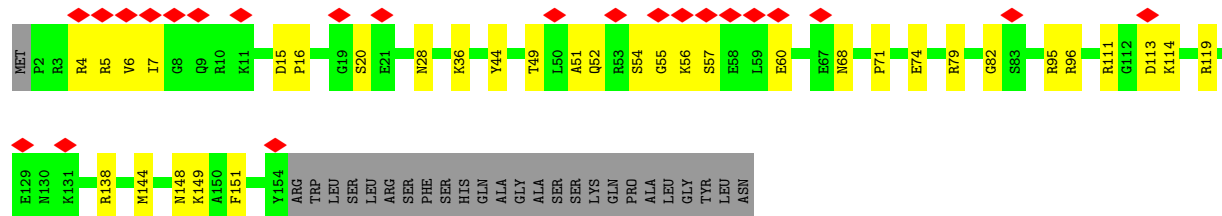
- Molecule 5: 30S ribosomal protein S5



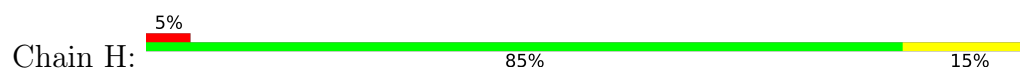
- Molecule 6: 30S ribosomal protein S6, fully modified isoform

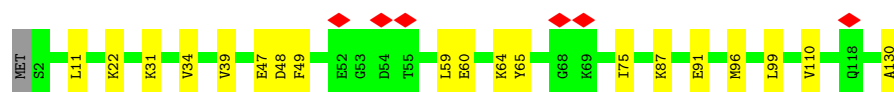


- Molecule 7: 30S ribosomal protein S7

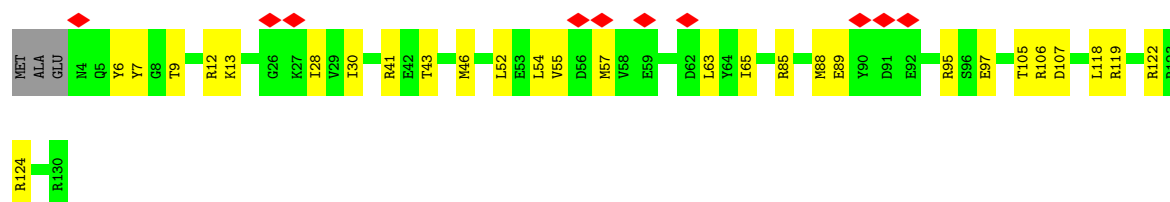
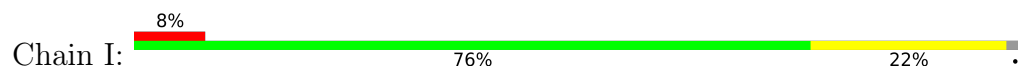


- Molecule 8: 30S ribosomal protein S8

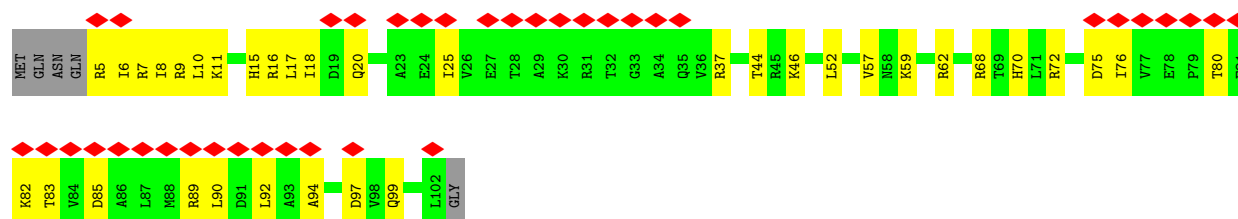




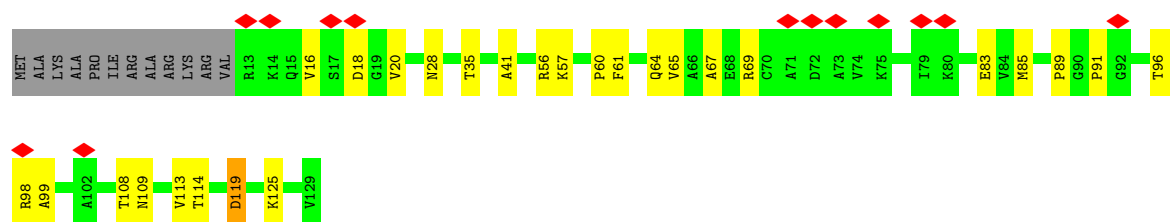
- Molecule 9: 30S ribosomal protein S9



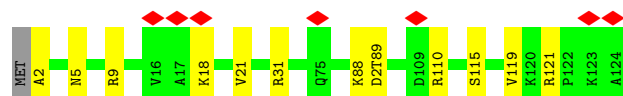
- Molecule 10: 30S ribosomal protein S10



- Molecule 11: 30S ribosomal protein S11

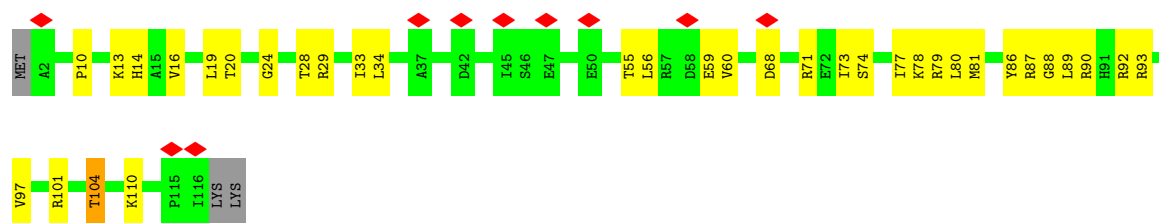


- Molecule 12: 30S ribosomal protein S12

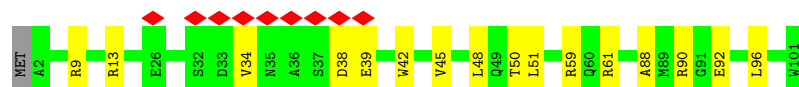
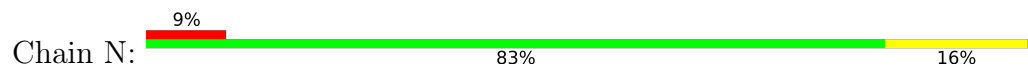


- Molecule 13: 30S ribosomal protein S13

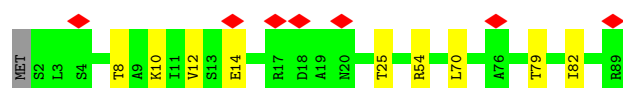
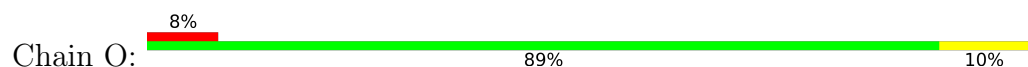




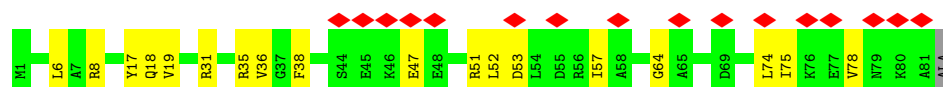
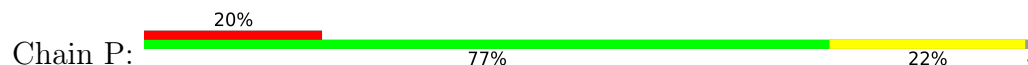
- Molecule 14: 30S ribosomal protein S14



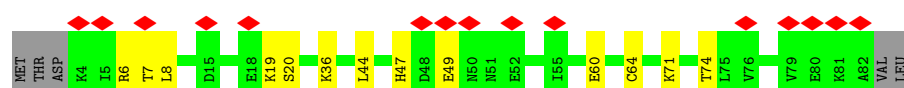
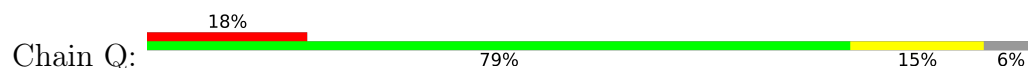
- Molecule 15: 30S ribosomal protein S15



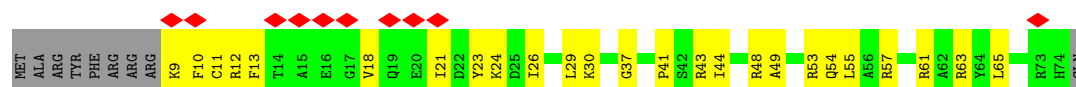
- Molecule 16: 30S ribosomal protein S16



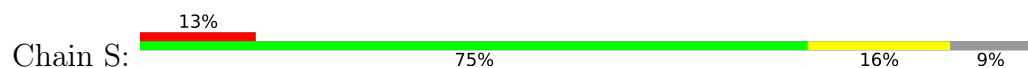
- Molecule 17: 30S ribosomal protein S17

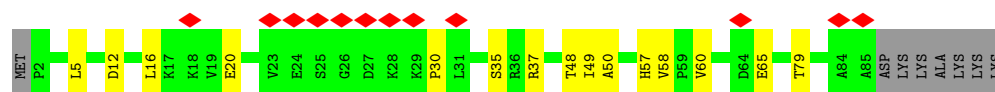


- Molecule 18: 30S ribosomal protein S18

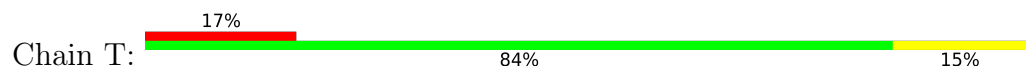


- Molecule 19: 30S ribosomal protein S19

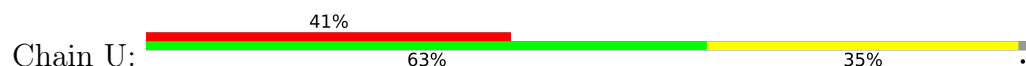




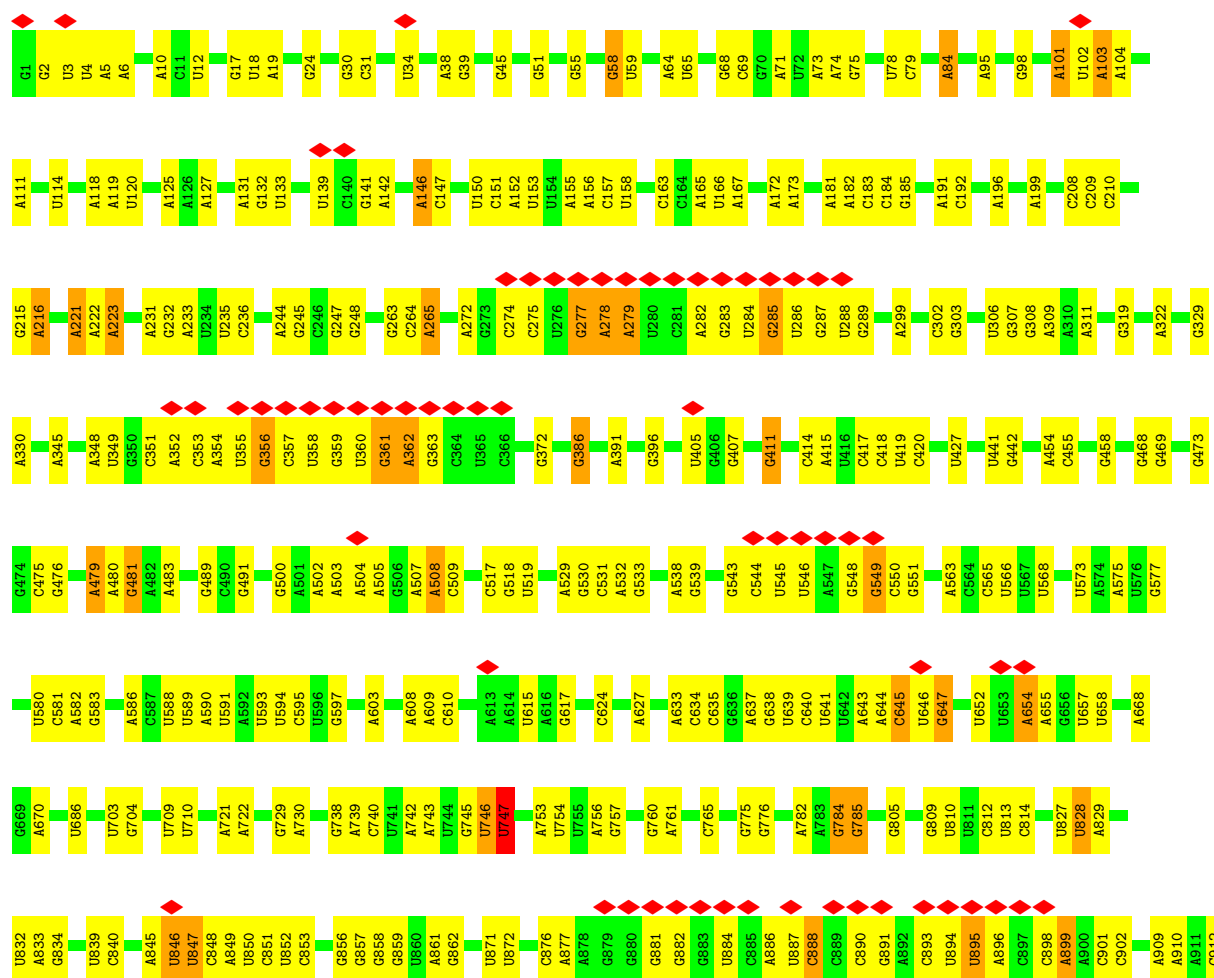
- Molecule 20: 30S ribosomal protein S20



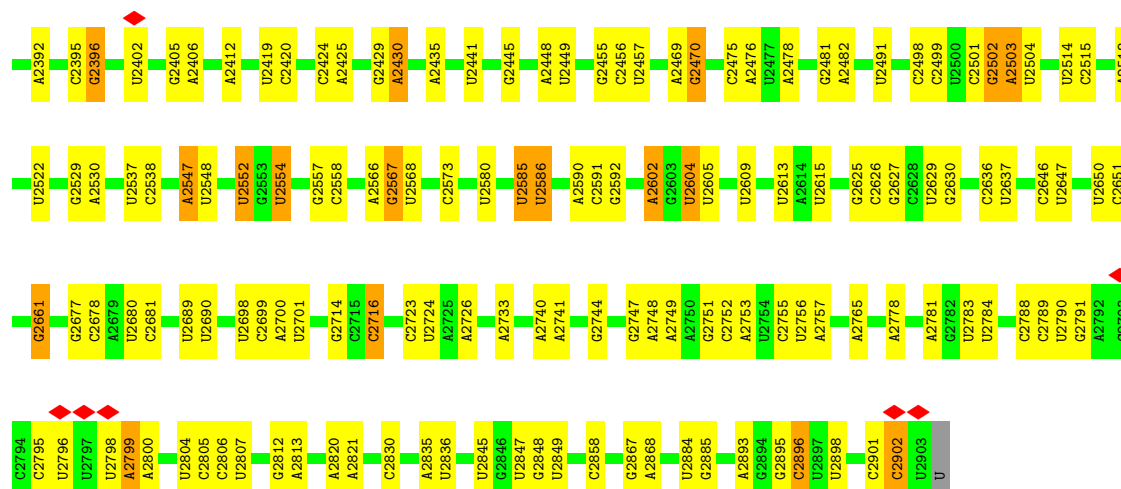
- Molecule 21: 30S ribosomal protein S21



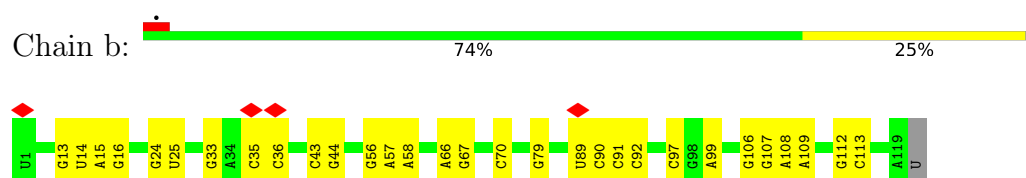
- Molecule 22: 23S rRNA



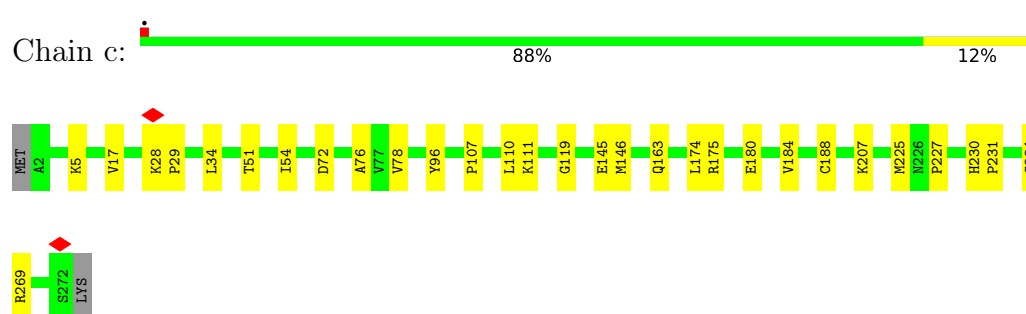




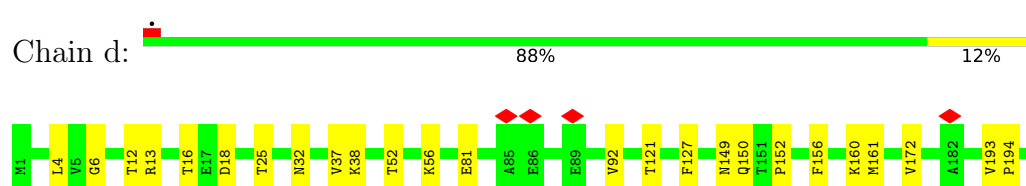
• Molecule 23: 5S rRNA



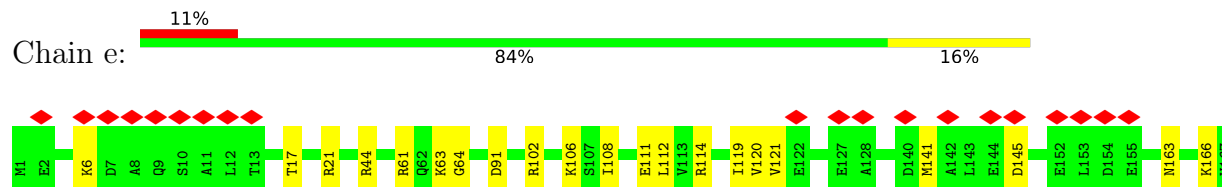
• Molecule 24: 50S ribosomal protein L2



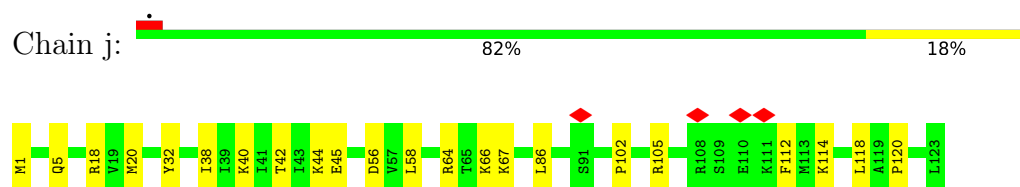
• Molecule 25: 50S ribosomal protein L3



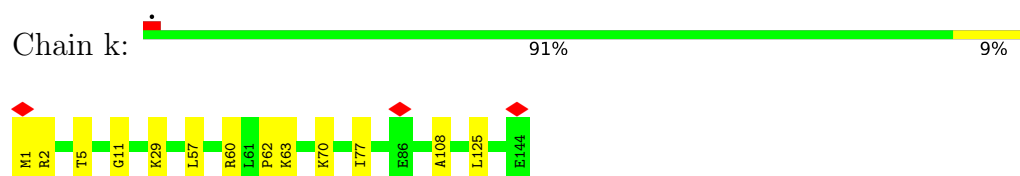
• Molecule 26: 50S ribosomal protein L4



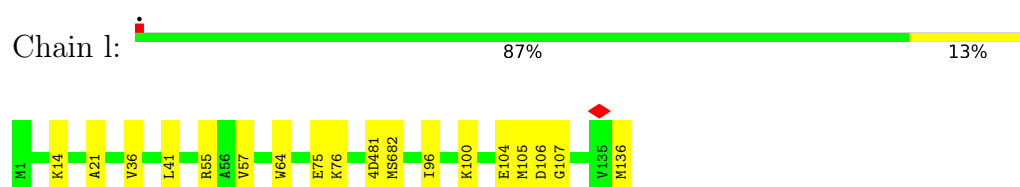
- Molecule 31: 50S ribosomal protein L14



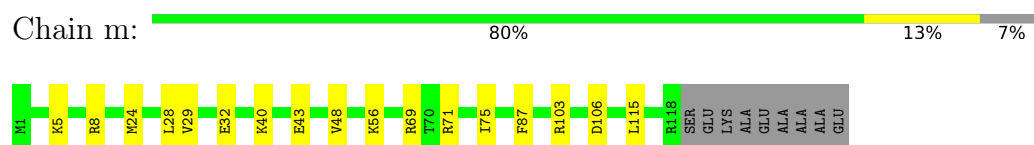
- Molecule 32: 50S ribosomal protein L15



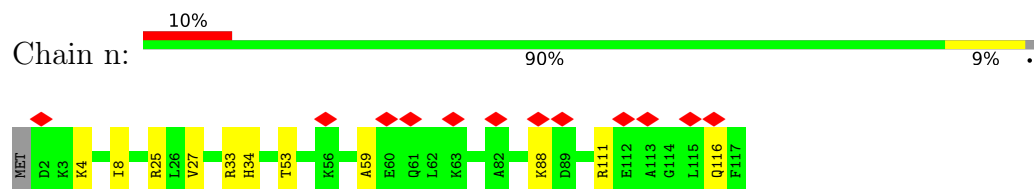
- Molecule 33: 50S ribosomal protein L16



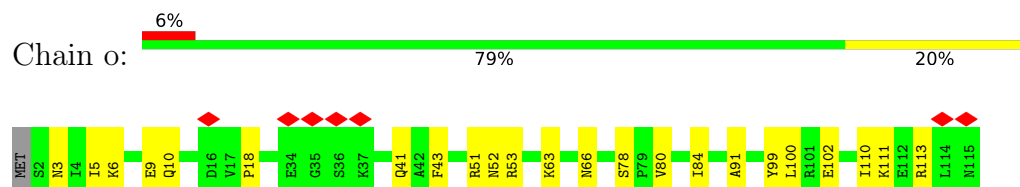
- Molecule 34: 50S ribosomal protein L17



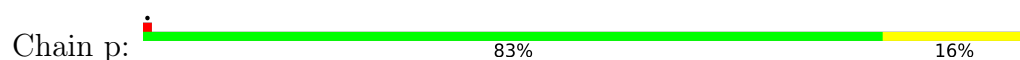
- Molecule 35: 50S ribosomal protein L18



- Molecule 36: 50S ribosomal protein L19

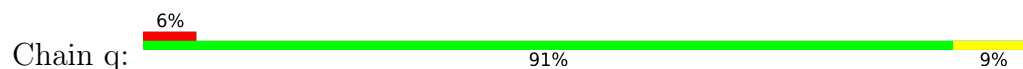


- Molecule 37: 50S ribosomal protein L20

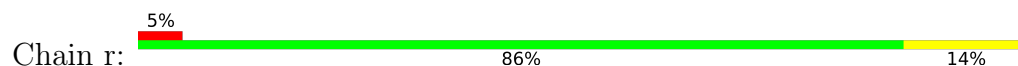




- Molecule 38: 50S ribosomal protein L21



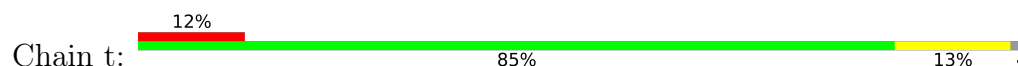
- Molecule 39: 50S ribosomal protein L22



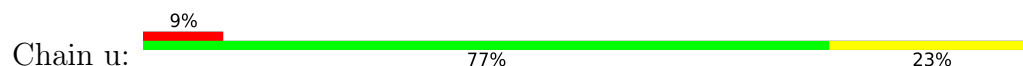
- Molecule 40: 50S ribosomal protein L23



- Molecule 41: 50S ribosomal protein L24



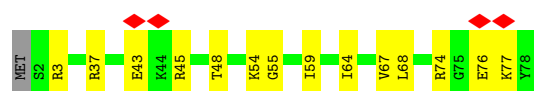
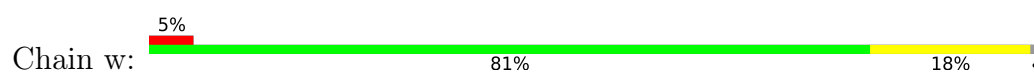
- Molecule 42: 50S ribosomal protein L25



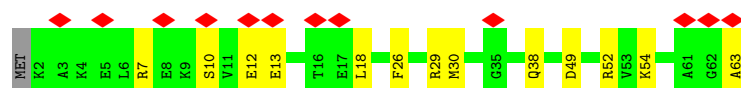
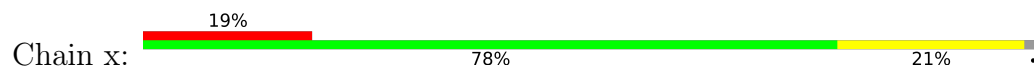
- Molecule 43: 50S ribosomal protein L27



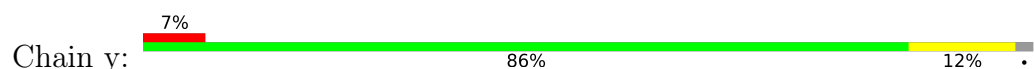
- Molecule 44: 50S ribosomal protein L28



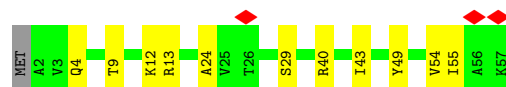
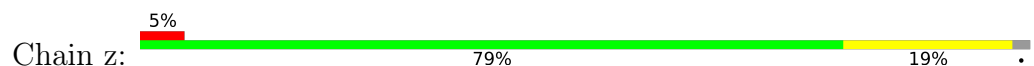
- Molecule 45: 50S ribosomal protein L29



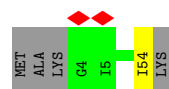
- Molecule 46: 50S ribosomal protein L30



- Molecule 47: 50S ribosomal protein L32



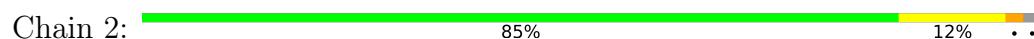
- Molecule 48: 50S ribosomal protein L33



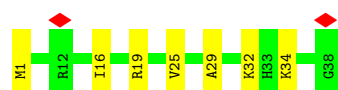
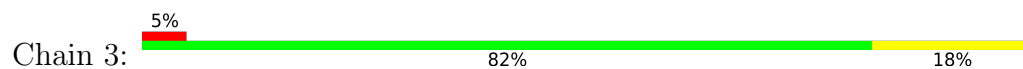
- Molecule 49: 50S ribosomal protein L34



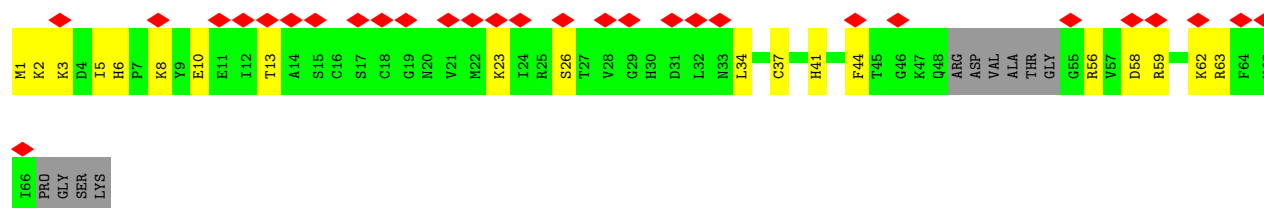
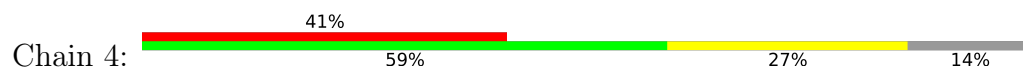
- Molecule 50: 50S ribosomal protein L35



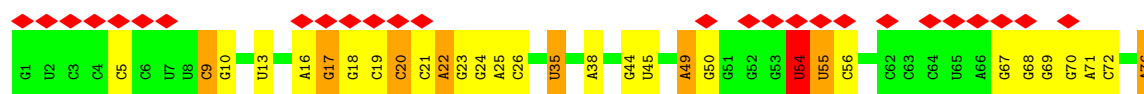
- Molecule 51: 50S ribosomal protein L36



- Molecule 52: 50S ribosomal protein L31



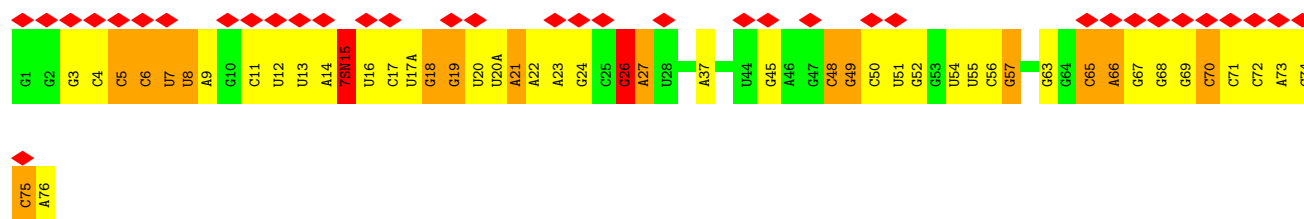
- Molecule 53: P-site tRNA_Glu



- Molecule 54: mRNA



- Molecule 55: A-site tRNA_Ile2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	260160	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.265	Depositor
Minimum map value	-0.138	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.0253	Depositor
Map size (Å)	439.10498, 439.10498, 439.10498	wwPDB
Map dimensions	530, 530, 530	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8285, 0.8285, 0.8285	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: T6A, IAS, U8U, 1MG, 4OC, OMG, G7M, MS6, 3TD, 7S3, OMU, 2MG, OMC, PSU, B8H, H2U, 2MA, 7SN, K, 4D4, MEQ, D2T, 6MZ, MG, UR3, MA6, 5MC, 5MU, OBX, M2G

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	A	0.19	0/36236	0.32	0/56520
2	B	0.15	0/1784	0.36	0/2403
3	C	0.17	0/1651	0.33	0/2225
4	D	0.14	0/1665	0.33	0/2227
5	E	0.18	0/1165	0.37	0/1568
6	F	0.16	0/858	0.36	0/1160
7	G	0.17	0/1219	0.36	0/1635
8	H	0.16	0/989	0.37	0/1326
9	I	0.17	0/1034	0.40	0/1375
10	J	0.18	0/796	0.43	0/1077
11	K	0.15	0/884	0.35	0/1191
12	L	0.16	0/960	0.38	0/1286
13	M	0.19	0/900	0.39	0/1204
14	N	0.19	0/817	0.37	0/1088
15	O	0.19	0/722	0.39	0/964
16	P	0.18	0/653	0.37	0/877
17	Q	0.14	0/650	0.37	0/871
18	R	0.16	0/553	0.34	0/742
19	S	0.16	0/685	0.33	0/922
20	T	0.18	0/676	0.34	0/895
21	U	0.16	0/597	0.35	0/792
22	a	0.22	1/65651 (0.0%)	0.34	0/102413
23	b	0.18	0/2850	0.29	0/4444
24	c	0.18	0/2121	0.38	0/2852
25	d	0.20	0/1576	0.38	0/2119
26	e	0.17	0/1571	0.34	0/2113
27	f	0.17	0/1434	0.34	0/1926
28	g	0.16	0/1343	0.42	0/1816
29	h	0.18	0/306	0.45	0/413
30	i	0.18	0/1152	0.35	0/1551
31	j	0.20	0/955	0.41	0/1279

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	k	0.20	0/1062	0.40	0/1413
33	l	0.18	0/1073	0.36	0/1433
34	m	0.19	0/958	0.39	0/1281
35	n	0.16	0/902	0.36	0/1209
36	o	0.17	0/929	0.33	0/1242
37	p	0.20	0/960	0.35	0/1278
38	q	0.17	0/829	0.39	0/1107
39	r	0.21	0/864	0.38	0/1156
40	s	0.17	0/744	0.38	0/994
41	t	0.18	0/787	0.47	0/1051
42	u	0.16	0/766	0.34	0/1025
43	v	0.17	0/593	0.32	0/785
44	w	0.17	0/635	0.32	0/848
45	x	0.17	0/502	0.35	0/667
46	y	0.20	0/453	0.42	0/605
47	z	0.19	0/450	0.36	0/599
48	0	0.16	0/424	0.32	0/565
49	1	0.17	0/380	0.35	0/498
50	2	0.20	0/513	0.43	0/676
51	3	0.18	0/303	0.37	0/397
52	4	0.14	0/488	0.32	0/649
53	V	0.15	0/1686	0.30	0/2621
54	W	0.20	0/264	0.25	0/409
55	X	0.20	0/1639	0.36	0/2545
All	All	0.20	1/152657 (0.0%)	0.34	0/228327

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
13	M	0	1
28	g	0	3
50	2	0	1
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	2069	G7M	O3'-P	5.19	1.61	1.56

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
50	2	31	HIS	Peptide
2	B	86	SER	Peptide
13	M	104	THR	Peptide
28	g	46	ALA	Peptide
28	g	47	ASP	Peptide
28	g	48	ASN	Peptide

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	A	32612	0	16432	317	0
2	B	1753	0	1780	34	0
3	C	1624	0	1696	19	0
4	D	1643	0	1707	30	0
5	E	1152	0	1196	13	0
6	F	839	0	833	19	0
7	G	1203	0	1254	25	0
8	H	979	0	1031	12	0
9	I	1022	0	1070	21	0
10	J	786	0	828	24	0
11	K	877	0	884	21	0
12	L	957	0	1017	11	0
13	M	891	0	952	25	0
14	N	805	0	844	11	0
15	O	714	0	734	5	0
16	P	643	0	661	14	0
17	Q	641	0	682	11	0
18	R	544	0	565	15	0
19	S	668	0	693	11	0
20	T	670	0	719	9	0
21	U	589	0	629	20	0
22	a	59130	0	29767	480	0
23	b	2549	0	1291	12	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
24	c	2082	0	2154	21	0
25	d	1566	0	1618	17	0
26	e	1552	0	1619	24	0
27	f	1410	0	1444	32	0
28	g	1323	0	1371	33	0
29	h	303	0	327	3	0
30	i	1129	0	1162	16	0
31	j	946	0	1023	16	0
32	k	1053	0	1129	14	0
33	l	1075	0	1145	11	0
34	m	945	0	989	11	0
35	n	892	0	923	6	0
36	o	917	0	962	15	0
37	p	947	0	1019	16	0
38	q	816	0	839	8	0
39	r	857	0	922	9	0
40	s	738	0	807	16	0
41	t	779	0	831	9	0
42	u	753	0	780	13	0
43	v	586	0	596	1	0
44	w	625	0	652	9	0
45	x	501	0	531	9	0
46	y	449	0	488	6	0
47	z	444	0	458	11	0
48	0	417	0	451	1	0
49	1	377	0	418	6	0
50	2	504	0	572	7	0
51	3	302	0	343	5	0
52	4	480	0	482	14	0
53	V	1619	0	832	15	0
54	W	236	0	119	2	0
55	X	1687	0	813	26	0
56	A	94	0	0	0	0
56	N	1	0	0	0	0
56	V	2	0	0	0	0
56	W	2	0	0	0	0
56	X	2	0	0	0	0
56	a	208	0	0	0	0
56	b	5	0	0	0	0
56	d	1	0	0	0	0
56	p	1	0	0	0	0
56	z	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
57	A	1	0	0	0	0
All	All	141919	0	95084	1384	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 6.

All (1384) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:G:H1	1:A:93:U:H3	1.12	0.95
53:V:35:U8U:HN3	54:W:30:G:H22	1.07	0.93
22:a:881:G:H1	22:a:895:U:H3	0.96	0.92
2:B:18:HIS:ND1	2:B:188:ASP:OD2	2.08	0.85
22:a:882:G:H1	22:a:894:U:H3	1.27	0.83
2:B:111:ILE:HG12	2:B:148:LEU:HD21	1.61	0.83
1:A:407:U:O2'	4:D:113:GLU:OE1	1.97	0.82
1:A:1441:A:H62	1:A:1461:G:H21	1.27	0.81
32:k:63:LYS:HE2	50:2:12:LYS:HD3	1.63	0.79
1:A:673:A:H2'	1:A:674:G:C8	2.18	0.78
24:c:78:VAL:HG21	24:c:110:LEU:HD21	1.65	0.78
9:I:106:ARG:NH1	9:I:107:ASP:O	2.16	0.78
22:a:279:A:OP2	22:a:361:G:N2	2.16	0.77
34:m:32:GLU:HG2	34:m:115:LEU:HD12	1.68	0.76
22:a:881:G:O6	22:a:895:U:O4	2.04	0.75
1:A:1009:U:H3	1:A:1020:G:H1	1.35	0.75
9:I:52:LEU:HD11	9:I:63:LEU:HD11	1.67	0.75
30:i:31:GLU:OE2	30:i:35:ARG:NE	2.16	0.75
17:Q:47:HIS:HB3	17:Q:74:THR:HG22	1.69	0.75
1:A:1086:U:H3	1:A:1099:G:H22	1.33	0.74
9:I:54:LEU:HD21	9:I:97:GLU:HG3	1.68	0.74
44:w:74:ARG:NH2	44:w:76:GLU:OE2	2.20	0.74
1:A:677:U:H3	1:A:713:G:H22	1.33	0.74
1:A:1097:C:H5''	2:B:139:ARG:HH22	1.53	0.73
7:G:144:MET:O	7:G:148:ASN:ND2	2.21	0.73
4:D:172:GLU:HG3	4:D:183:LYS:HZ2	1.54	0.73
17:Q:7:THR:HG23	17:Q:60:GLU:HG3	1.71	0.73
22:a:1168:G:H1	22:a:1181:U:H3	1.34	0.72
13:M:101:ARG:NE	13:M:104:THR:OG1	2.19	0.72
31:j:114:LYS:HE2	31:j:118:LEU:HD11	1.72	0.72
19:S:5:LEU:HD21	52:4:63:ARG:HH21	1.55	0.72
1:A:202:G:O2'	1:A:468:A:N3	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:1607:C:N4	22:a:1622:G:OP2	2.22	0.72
42:u:9:ARG:HG2	42:u:41:GLU:HG3	1.72	0.72
10:J:80:THR:HG22	10:J:82:LYS:H	1.55	0.71
28:g:107:LEU:O	28:g:152:ARG:NH1	2.23	0.71
1:A:1119:C:OP1	9:I:85:ARG:NH1	2.23	0.71
55:X:51:U:H3	55:X:63:G:H1	0.77	0.71
1:A:736:C:OP1	18:R:61:ARG:NH1	2.24	0.71
2:B:86:SER:O	2:B:222:ARG:NH2	2.23	0.71
1:A:310:G:H5'	16:P:31:ARG:HB3	1.72	0.70
11:K:97:ILE:HG22	21:U:12:PHE:HZ	1.56	0.70
27:f:57:LEU:HD22	27:f:65:PRO:HB3	1.73	0.70
22:a:1153:C:OP1	37:p:92:ARG:NH2	2.25	0.69
3:C:77:ILE:HA	3:C:84:VAL:HG23	1.74	0.69
22:a:1869:G:N2	22:a:1872:A:OP2	2.22	0.69
45:x:49:ASP:OD1	45:x:52:ARG:NH1	2.25	0.69
1:A:664:G:H22	1:A:741:G:H1	1.41	0.69
4:D:125:VAL:HG12	4:D:143:VAL:HG12	1.75	0.69
8:H:47:GLU:HG3	8:H:64:LYS:HB3	1.75	0.69
22:a:475:C:O2	22:a:479:A:N6	2.26	0.68
22:a:287:G:O6	22:a:353:C:N4	2.16	0.68
26:e:119:ILE:HB	26:e:187:VAL:HG12	1.75	0.68
55:X:65:C:H2'	55:X:66:A:H8	1.59	0.68
8:H:96:MET:HG3	8:H:99:LEU:HB2	1.75	0.68
26:e:61:ARG:NH1	26:e:63:LYS:O	2.26	0.68
55:X:51:U:O4	55:X:63:G:O6	2.12	0.67
1:A:1124:G:OP1	10:J:37:ARG:NH2	2.28	0.67
30:i:125:TYR:OH	30:i:132:HIS:NE2	2.25	0.67
22:a:2547:A:H2'	22:a:2548:U:C6	2.29	0.67
22:a:1434:A:H2'	22:a:1435:G:H8	1.60	0.67
22:a:285:G:O6	22:a:355:U:O2	2.13	0.66
1:A:1391:U:H2'	1:A:1392:G:C8	2.29	0.66
22:a:287:G:N1	22:a:353:C:N3	2.34	0.66
1:A:545:C:OP1	4:D:58:LYS:NZ	2.29	0.66
1:A:626:G:OP1	16:P:18:GLN:NE2	2.29	0.66
1:A:946:A:H2'	1:A:947:G:C8	2.30	0.66
26:e:21:ARG:HH21	26:e:106:LYS:HD3	1.61	0.66
1:A:957:U:H4'	19:S:79:THR:HG23	1.78	0.66
13:M:68:ASP:OD1	13:M:71:ARG:NH2	2.28	0.66
22:a:153:U:OP1	44:w:77:LYS:NZ	2.21	0.66
22:a:2591:C:H2'	22:a:2592:G:C8	2.30	0.66
28:g:164:TYR:HB2	28:g:167:GLU:HB2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:A:H2'	1:A:460:A:C8	2.31	0.66
25:d:121:THR:HG22	25:d:161:MET:HA	1.78	0.66
1:A:73:C:HO2'	1:A:74:A:H8	1.42	0.65
2:B:117:LEU:HG	2:B:141:LEU:HD12	1.79	0.65
26:e:181:ILE:HG12	32:k:1:MET:HE2	1.79	0.65
13:M:71:ARG:NH1	27:f:113:ASP:OD2	2.29	0.65
47:z:54:VAL:HG23	47:z:55:ILE:HG12	1.78	0.65
1:A:1356:G:H2'	1:A:1357:A:C8	2.32	0.65
3:C:164:ARG:NE	3:C:166:GLU:OE2	2.28	0.65
6:F:41:ASP:OD1	6:F:58:HIS:NE2	2.28	0.65
22:a:652:U:OP1	22:a:654:A:N6	2.29	0.65
1:A:147:G:H2'	1:A:148:G:C8	2.32	0.65
21:U:63:GLU:HG3	21:U:66:ARG:HH21	1.61	0.65
29:h:1:MET:N	29:h:21:VAL:O	2.29	0.65
11:K:35:THR:HG22	11:K:41:ALA:HA	1.77	0.65
22:a:2830:C:H5''	25:d:56:LYS:HE3	1.78	0.64
22:a:210:C:OP1	49:l:29:GLN:NE2	2.30	0.64
22:a:1434:A:H2'	22:a:1435:G:C8	2.32	0.64
14:N:90:ARG:NH1	14:N:92:GLU:OE1	2.31	0.64
22:a:846:U:H4'	22:a:847:U:H5	1.62	0.64
26:e:145:ASP:HB3	26:e:166:LYS:HE2	1.78	0.64
55:X:15:7SN:N71	55:X:16:U:O4	2.30	0.64
52:4:59:ARG:HA	52:4:62:LYS:HE3	1.78	0.64
1:A:127:G:O2'	17:Q:6:ARG:NH1	2.30	0.64
22:a:84:A:N1	22:a:98:G:O2'	2.31	0.64
1:A:17:U:H2'	1:A:18:C:C6	2.34	0.64
24:c:29:PRO:HG2	24:c:34:LEU:HD11	1.79	0.63
1:A:437:U:HO2'	4:D:120:HIS:HD1	1.41	0.63
22:a:2000:C:OP1	34:m:5:LYS:NZ	2.30	0.63
1:A:1343:G:OP1	9:I:124:ARG:NH2	2.30	0.63
11:K:109:ASN:HD21	21:U:5:LYS:HG3	1.64	0.63
26:e:119:ILE:HD11	26:e:185:LYS:HD3	1.80	0.63
28:g:16:ASP:HB3	28:g:27:LYS:HB3	1.81	0.63
6:F:38:ARG:HG2	6:F:63:ASN:HB2	1.82	0.62
22:a:568:U:H1'	22:a:2030:6MZ:H9C1	1.81	0.62
22:a:1469:A:H2'	22:a:1470:A:C8	2.34	0.62
1:A:28:A:O2'	1:A:296:U:OP1	2.17	0.62
22:a:282:A:H2'	22:a:283:G:C8	2.35	0.62
23:b:79:G:N7	42:u:14:LYS:NZ	2.47	0.62
2:B:188:ASP:CG	2:B:189:THR:H	2.07	0.62
4:D:198:HIS:O	4:D:202:GLU:HG3	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:662:U:O2'	1:A:836:G:OP1	2.17	0.62
9:I:7:TYR:N	9:I:89:GLU:OE1	2.33	0.62
53:V:69:G:H2'	53:V:70:G:H8	1.65	0.62
55:X:65:C:H2'	55:X:66:A:C8	2.35	0.62
45:x:18:LEU:HD11	45:x:54:LYS:HE3	1.81	0.61
22:a:172:A:H2'	22:a:173:A:C8	2.36	0.61
22:a:483:A:C8	41:t:45:HIS:HD2	2.19	0.61
3:C:135:LYS:O	3:C:139:GLN:HG2	1.99	0.61
22:a:2313:C:H5''	27:f:88:LYS:HD3	1.81	0.61
45:x:29:ARG:HG2	45:x:30:MET:HE2	1.83	0.61
11:K:98:ARG:NH1	21:U:13:ASP:OD1	2.33	0.61
28:g:52:PHE:CZ	28:g:72:LEU:HD22	2.36	0.61
1:A:150:U:H2'	1:A:151:A:H8	1.65	0.61
1:A:337:G:H2'	1:A:338:A:C8	2.36	0.61
31:j:18:ARG:HG2	31:j:45:GLU:OE1	2.00	0.61
26:e:6:LYS:HG2	26:e:121:VAL:HG12	1.83	0.61
34:m:28:LEU:HD23	34:m:48:VAL:HG21	1.83	0.61
22:a:639:U:H2'	22:a:640:C:H6	1.66	0.60
1:A:197:A:N1	1:A:220:G:O2'	2.31	0.60
1:A:235:C:H2'	1:A:236:A:C8	2.36	0.60
22:a:1141:U:H4'	22:a:1142:A:O4'	2.01	0.60
1:A:131:A:H2'	1:A:132:C:C6	2.36	0.60
22:a:742:A:H2'	22:a:743:A:C8	2.36	0.60
22:a:1720:U:H2'	22:a:1721:G:O4'	2.02	0.60
25:d:12:THR:OG1	25:d:13:ARG:N	2.35	0.60
28:g:22:GLN:NE2	28:g:38:ASN:O	2.34	0.60
1:A:154:U:H2'	1:A:155:A:C8	2.37	0.60
22:a:1536:C:H4'	22:a:1537:G:H5''	1.84	0.60
1:A:167:A:N6	1:A:168:G:O6	2.35	0.60
22:a:274:C:H2'	22:a:275:C:C6	2.36	0.60
22:a:2469:A:H4'	33:l:55:ARG:NE	2.17	0.60
3:C:50:ALA:HA	3:C:75:ILE:HD11	1.83	0.60
1:A:127:G:H4'	17:Q:6:ARG:HH12	1.67	0.60
30:i:13:ARG:NH2	30:i:49:ASP:O	2.35	0.60
30:i:23:LYS:NZ	30:i:142:ILE:OXT	2.30	0.60
22:a:848:C:H2'	22:a:849:A:H8	1.68	0.59
1:A:1323:G:H2'	1:A:1324:A:C8	2.37	0.59
10:J:5:ARG:HH22	10:J:7:ARG:HH21	1.48	0.59
28:g:18:LYS:HB3	28:g:25:THR:HB	1.84	0.59
22:a:1802:A:H2'	22:a:1803:A:C8	2.37	0.59
1:A:363:A:OP2	12:L:31:ARG:NH1	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:b:66:A:N6	23:b:107:G:H2'	2.18	0.59
27:f:111:ILE:HB	27:f:114:PHE:HB2	1.82	0.59
1:A:56:U:H2'	1:A:57:G:C8	2.38	0.59
1:A:1009:U:O2	1:A:1020:G:N2	2.34	0.59
25:d:16:THR:OG1	25:d:18:ASP:OD1	2.20	0.59
3:C:168:TYR:OH	5:E:55:GLU:OE2	2.17	0.59
22:a:483:A:H5''	41:t:47:LYS:HD2	1.85	0.59
22:a:2328:A:H2'	22:a:2329:U:C6	2.37	0.59
22:a:191:A:H2'	22:a:192:C:C6	2.36	0.59
10:J:37:ARG:HB2	10:J:75:ASP:HB2	1.85	0.59
1:A:501:C:H2'	1:A:502:A:H8	1.68	0.59
1:A:714:G:H2'	1:A:715:A:C8	2.37	0.59
11:K:108:THR:O	11:K:109:ASN:ND2	2.36	0.59
15:O:8:THR:O	15:O:12:VAL:HG23	2.03	0.59
1:A:80:A:H2'	1:A:81:A:C8	2.38	0.58
1:A:477:C:H2'	1:A:478:A:C8	2.38	0.58
1:A:1106:G:H5''	3:C:172:ARG:HG2	1.85	0.58
22:a:299:A:N3	22:a:319:G:O2'	2.35	0.58
27:f:147:ASP:OD1	27:f:148:ARG:N	2.36	0.58
40:s:56:GLU:HB3	40:s:86:THR:HG23	1.83	0.58
55:X:6:C:H2'	55:X:7:U:H6	1.68	0.58
1:A:519:C:O2'	1:A:520:A:OP1	2.21	0.58
1:A:544:G:OP1	4:D:56:ARG:NH2	2.35	0.58
1:A:1397:C:H41	54:W:35:A:H5''	1.67	0.58
1:A:981:U:OP1	14:N:9:ARG:NH1	2.36	0.58
24:c:163:GLN:OE1	24:c:175:ARG:NH1	2.35	0.58
1:A:1151:A:H5''	10:J:44:THR:HG23	1.85	0.58
7:G:16:PRO:HA	9:I:46:MET:SD	2.43	0.58
22:a:2306:C:OP2	22:a:2307:G:O2'	2.20	0.58
27:f:62:GLY:O	27:f:95:ARG:NH1	2.36	0.58
1:A:728:A:H2'	1:A:729:A:C8	2.39	0.58
1:A:1008:U:H3'	1:A:1009:U:H5''	1.85	0.58
1:A:1031:C:O3'	1:A:1032:G:N2	2.36	0.58
1:A:1033:G:H2'	1:A:1034:G:O4'	2.02	0.58
22:a:1013:C:H2'	22:a:1014:A:H8	1.68	0.58
22:a:1797:G:HO2'	24:c:257:THR:HG1	1.50	0.58
33:l:21:ALA:HB1	33:l:100:LYS:HD3	1.84	0.58
10:J:16:ARG:O	10:J:20:GLN:HG3	2.03	0.58
22:a:593:U:H2'	22:a:594:U:C6	2.39	0.58
22:a:639:U:H2'	22:a:640:C:C6	2.38	0.58
26:e:17:THR:HB	26:e:21:ARG:NH2	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:b:66:A:H61	23:b:107:G:H2'	1.68	0.58
22:a:956:G:H5'	33:l:76:LYS:HD2	1.84	0.58
28:g:52:PHE:HZ	28:g:72:LEU:HD22	1.69	0.58
1:A:404:G:OP2	4:D:115:ARG:NH2	2.31	0.57
1:A:492:C:H2'	1:A:493:A:C8	2.39	0.57
22:a:1168:G:H2'	22:a:1169:A:C8	2.39	0.57
13:M:79:ARG:HD2	19:S:65:GLU:OE1	2.05	0.57
22:a:856:G:H2'	22:a:857:G:C8	2.40	0.57
22:a:2756:U:OP2	51:3:19:ARG:NH2	2.36	0.57
26:e:108:ILE:HG12	32:k:1:MET:HE1	1.85	0.57
32:k:2:ARG:H	32:k:5:THR:HG22	1.69	0.57
1:A:1352:C:H2'	1:A:1353:G:C8	2.40	0.57
22:a:2798:U:H4'	22:a:2799:A:H5'	1.86	0.57
25:d:25:THR:HG21	25:d:193:VAL:HG13	1.86	0.57
1:A:501:C:H2'	1:A:502:A:C8	2.38	0.57
1:A:501:C:OP2	12:L:121:ARG:NH1	2.38	0.57
1:A:1287:A:H2'	1:A:1288:A:C8	2.40	0.57
2:B:70:VAL:HG12	2:B:169:GLU:HG2	1.86	0.57
7:G:5:ARG:HH22	7:G:7:ILE:HB	1.69	0.57
10:J:85:ASP:HB3	10:J:89:ARG:HH22	1.69	0.57
10:J:52:LEU:HD21	10:J:59:LYS:HD2	1.86	0.57
41:t:14:LEU:HD11	41:t:71:ALA:HB2	1.86	0.57
50:2:62:LEU:HB3	50:2:65:ALA:HB3	1.86	0.57
1:A:56:U:H2'	1:A:57:G:H8	1.68	0.57
1:A:181:A:N6	1:A:195:A:OP2	2.38	0.57
22:a:1000:A:H2'	22:a:1001:A:C8	2.40	0.57
22:a:1570:A:H2'	22:a:1571:A:C8	2.40	0.57
31:j:38:ILE:HD11	31:j:112:PHE:HZ	1.69	0.57
55:X:65:C:O2'	55:X:66:A:OP1	2.22	0.57
55:X:6:C:H2'	55:X:7:U:C6	2.40	0.57
1:A:1124:G:N2	1:A:1125:U:O4	2.34	0.57
22:a:2071:A:H2'	22:a:2072:C:C6	2.39	0.57
55:X:51:U:O2	55:X:63:G:N2	2.29	0.57
22:a:876:C:H2'	22:a:877:A:O4'	2.05	0.56
22:a:1365:A:OP1	44:w:3:ARG:NH1	2.32	0.56
22:a:78:U:H5'	45:x:7:ARG:HH22	1.69	0.56
22:a:282:A:H2'	22:a:283:G:H8	1.70	0.56
22:a:2243:U:H2'	22:a:2244:U:C6	2.39	0.56
1:A:1218:C:H2'	1:A:1219:A:C8	2.40	0.56
13:M:33:ILE:HD13	13:M:60:VAL:HG22	1.87	0.56
22:a:2291:U:H2'	22:a:2292:U:C6	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:2627:G:O2'	22:a:2781:A:N1	2.33	0.56
28:g:149:ARG:HA	28:g:162:VAL:HB	1.88	0.56
36:o:51:ARG:HG2	36:o:53:ARG:HG2	1.88	0.56
6:F:40:GLU:OE1	6:F:100:SER:OG	2.21	0.56
9:I:6:TYR:HD2	9:I:89:GLU:HG2	1.70	0.56
22:a:155:A:H2'	22:a:156:A:C8	2.40	0.56
22:a:2537:U:H2'	22:a:2538:C:C6	2.40	0.56
1:A:910:C:OP2	12:L:18:LYS:NZ	2.31	0.56
15:O:25:THR:HG21	15:O:70:LEU:HB2	1.88	0.56
1:A:193:C:H4'	20:T:56:PRO:HG3	1.87	0.56
1:A:269:C:H2'	1:A:270:A:C8	2.41	0.56
22:a:1412:U:H2'	22:a:1413:A:C8	2.40	0.56
1:A:255:G:H4'	17:Q:19:LYS:HD3	1.87	0.56
28:g:41:VAL:O	28:g:55:ARG:NH1	2.39	0.56
22:a:989:G:OP2	46:y:12:SER:OG	2.16	0.56
1:A:31:G:O2'	1:A:48:C:N4	2.39	0.56
1:A:1273:C:H2'	1:A:1274:A:O4'	2.06	0.55
22:a:1224:U:H2'	22:a:1225:G:C8	2.41	0.55
22:a:2376:A:N3	35:n:111:ARG:NH2	2.53	0.55
24:c:225:MET:SD	24:c:230:HIS:HB2	2.45	0.55
1:A:73:C:O2'	1:A:74:A:H8	1.89	0.55
15:O:79:THR:O	15:O:82:ILE:HG22	2.05	0.55
22:a:634:C:H2'	22:a:635:C:C6	2.41	0.55
23:b:43:C:H5''	52:4:1:MET:HG2	1.87	0.55
5:E:13:GLU:HG2	5:E:39:VAL:HG22	1.88	0.55
22:a:1178:C:H2'	22:a:1179:G:H8	1.71	0.55
27:f:4:LEU:HD13	27:f:101:GLU:HB2	1.87	0.55
1:A:236:A:H2'	1:A:237:G:C8	2.41	0.55
9:I:43:THR:HG22	9:I:46:MET:HE2	1.88	0.55
22:a:2591:C:H2'	22:a:2592:G:H8	1.71	0.55
1:A:384:G:H2'	1:A:385:C:C6	2.41	0.55
1:A:461:A:H2'	1:A:462:G:H8	1.71	0.55
22:a:1168:G:H2'	22:a:1169:A:H8	1.72	0.55
1:A:391:G:H5''	16:P:8:ARG:HE	1.72	0.55
1:A:511:C:H4'	4:D:44:ARG:HH22	1.72	0.55
22:a:538:A:H5''	30:i:7:LYS:HE3	1.87	0.55
22:a:1534:U:H1'	22:a:1538:G:N2	2.21	0.55
1:A:1124:G:H1'	1:A:1125:U:H5	1.71	0.55
22:a:657:U:H2'	22:a:658:U:C6	2.41	0.55
40:s:24:MET:SD	40:s:30:ILE:HG22	2.47	0.55
22:a:2377:A:H2'	22:a:2378:A:C8	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:k:57:LEU:HD22	50:2:54:ASP:HB3	1.88	0.55
53:V:69:G:H2'	53:V:70:G:C8	2.42	0.55
1:A:475:C:H2'	1:A:476:U:C6	2.41	0.55
1:A:1010:U:H2'	1:A:1011:C:C6	2.40	0.55
47:z:43:ILE:HG22	47:z:49:TYR:HB2	1.88	0.55
55:X:6:C:HO2'	55:X:7:U:C5'	2.20	0.55
1:A:223:A:H2'	1:A:224:U:C6	2.43	0.55
4:D:170:TRP:CE2	4:D:186:PRO:HG3	2.42	0.54
22:a:1300:G:H4'	22:a:1301:A:H5'	1.87	0.54
27:f:29:PRO:HB2	27:f:169:LEU:HD22	1.87	0.54
1:A:945:G:C2	1:A:946:A:C8	2.95	0.54
11:K:109:ASN:ND2	21:U:5:LYS:HG3	2.22	0.54
16:P:18:GLN:HG2	16:P:35:ARG:HE	1.72	0.54
22:a:933:A:H5'	22:a:934:U:OP2	2.07	0.54
40:s:54:GLU:OE1	40:s:91:GLN:NE2	2.40	0.54
22:a:58:G:H2'	22:a:59:U:C6	2.43	0.54
7:G:51:ALA:O	7:G:55:GLY:N	2.39	0.54
22:a:608:A:H2'	22:a:609:A:C8	2.42	0.54
22:a:2602:A:N7	53:V:76:A:H4'	2.22	0.54
1:A:946:A:H2'	1:A:947:G:H8	1.73	0.54
28:g:104:ASN:ND2	28:g:114:ASP:OD1	2.36	0.54
1:A:470:C:H2'	1:A:471:U:C6	2.43	0.54
1:A:1348:U:H4'	9:I:122:ARG:HD2	1.89	0.54
1:A:1447:A:OP1	1:A:1448:C:N4	2.30	0.54
22:a:617:G:OP1	26:e:102:ARG:NE	2.28	0.54
22:a:633:A:H5''	32:k:70:LYS:HD2	1.89	0.54
22:a:871:U:H2'	22:a:872:U:C6	2.43	0.54
28:g:170:ARG:NH2	51:3:29:ALA:O	2.40	0.54
1:A:537:G:H2'	1:A:538:G:C8	2.43	0.54
1:A:1175:G:H2'	1:A:1176:A:H8	1.72	0.54
21:U:41:PRO:HA	21:U:44:GLU:HG3	1.89	0.54
22:a:2074:U:H2'	22:a:2075:U:C6	2.43	0.54
24:c:180:GLU:HB2	24:c:269:ARG:O	2.06	0.54
42:u:58:SER:O	42:u:73:LYS:NZ	2.40	0.54
1:A:674:G:H2'	1:A:675:A:H8	1.71	0.54
1:A:1410:A:H2'	1:A:1411:C:C6	2.43	0.54
2:B:23:TRP:CZ3	2:B:25:PRO:HA	2.42	0.54
10:J:46:LYS:NZ	10:J:68:ARG:HH21	2.06	0.54
22:a:306:U:H2'	22:a:307:G:O4'	2.07	0.54
36:o:5:ILE:O	36:o:9:GLU:HG3	2.07	0.54
18:R:41:PRO:HD2	18:R:44:ILE:HD12	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:G:H2'	1:A:300:A:C8	2.43	0.53
22:a:1046:A:H3'	22:a:1047:G:H5'	1.90	0.53
22:a:1413:A:H2'	22:a:1414:C:O4'	2.08	0.53
1:A:1062:U:H2'	1:A:1063:C:C6	2.43	0.53
1:A:216:U:H2'	1:A:217:C:C6	2.43	0.53
1:A:555:U:H2'	1:A:556:C:C6	2.43	0.53
7:G:68:ASN:O	7:G:138:ARG:NH1	2.41	0.53
22:a:1410:G:H2'	22:a:1411:U:C6	2.43	0.53
52:4:56:ARG:O	52:4:59:ARG:HG2	2.08	0.53
1:A:87:C:H2'	1:A:88:U:C5	2.43	0.53
22:a:2095:A:OP1	29:h:11:ASN:ND2	2.42	0.53
22:a:2329:U:H2'	22:a:2330:G:C8	2.43	0.53
42:u:4:ILE:HB	42:u:63:ILE:HG12	1.91	0.53
22:a:351:C:H2'	22:a:352:A:C8	2.43	0.53
22:a:2514:U:H2'	22:a:2515:C:C6	2.44	0.53
24:c:110:LEU:HD23	24:c:111:LYS:N	2.23	0.53
1:A:459:A:H2'	1:A:460:A:H8	1.69	0.53
8:H:96:MET:HE3	8:H:130:ALA:HB1	1.90	0.53
22:a:832:U:H2'	22:a:833:A:C8	2.44	0.53
22:a:998:C:P	37:p:92:ARG:HH12	2.32	0.53
22:a:1028:A:H2'	22:a:1029:A:C8	2.44	0.53
27:f:167:ARG:NH1	27:f:171:ALA:HB2	2.23	0.53
1:A:1026:G:O2'	1:A:1027:C:OP1	2.25	0.53
3:C:130:PHE:O	3:C:134:MET:HG3	2.09	0.53
42:u:20:LEU:HD21	42:u:41:GLU:HG2	1.90	0.53
4:D:170:TRP:NE1	4:D:186:PRO:HG3	2.23	0.53
7:G:6:VAL:HG12	7:G:6:VAL:O	2.09	0.53
31:j:5:GLN:HA	31:j:20:MET:HE3	1.91	0.53
1:A:155:A:H2'	1:A:156:C:C6	2.44	0.53
7:G:151:PHE:CZ	11:K:56:ARG:HG3	2.44	0.53
12:L:5:ASN:OD1	17:Q:36:LYS:NZ	2.42	0.53
21:U:13:ASP:O	21:U:17:ARG:HG2	2.08	0.53
22:a:538:A:H4'	30:i:7:LYS:HG3	1.91	0.53
34:m:24:MET:HE1	34:m:40:LYS:HD3	1.90	0.53
1:A:195:A:N3	1:A:222:C:O2'	2.39	0.53
1:A:475:C:H2'	1:A:476:U:H6	1.74	0.53
22:a:549:G:H2'	22:a:550:C:C6	2.44	0.53
33:l:57:VAL:HG21	33:l:105:MET:HE1	1.90	0.53
53:V:24:G:H2'	53:V:25:A:C8	2.44	0.53
4:D:19:LEU:HB2	4:D:21:LEU:HG	1.89	0.52
22:a:2:G:H2'	22:a:3:U:C6	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:947:A:H2'	22:a:948:C:C6	2.44	0.52
36:o:43:PHE:CE2	36:o:63:LYS:HE3	2.44	0.52
37:p:11:ARG:O	37:p:15:LYS:HG3	2.09	0.52
1:A:537:G:H2'	1:A:538:G:H8	1.74	0.52
1:A:1180:A:OP1	9:I:105:THR:OG1	2.16	0.52
22:a:286:U:H2'	22:a:287:G:H8	1.75	0.52
27:f:16:LEU:HD12	27:f:28:VAL:HG13	1.90	0.52
1:A:568:G:O6	12:L:2:ALA:N	2.43	0.52
1:A:1524:C:H2'	1:A:1525:G:C8	2.45	0.52
17:Q:64:CYS:SG	17:Q:74:THR:HG23	2.49	0.52
51:3:16:ILE:HD13	51:3:25:VAL:HG22	1.90	0.52
5:E:88:VAL:HG22	5:E:93:ARG:HG2	1.91	0.52
22:a:1541:C:H2'	22:a:1542:U:C6	2.44	0.52
22:a:1794:A:H2'	22:a:1795:C:C6	2.45	0.52
27:f:3:LYS:HD2	27:f:4:LEU:HD12	1.91	0.52
1:A:399:G:H2'	1:A:400:C:C6	2.44	0.52
1:A:666:G:H5'	1:A:726:C:H1'	1.90	0.52
6:F:38:ARG:NH1	6:F:100:SER:HB3	2.24	0.52
22:a:909:A:H2'	22:a:912:C:H5	1.74	0.52
22:a:1754:A:N1	22:a:2716:C:O2'	2.39	0.52
55:X:11:C:H2'	55:X:12:U:H6	1.75	0.52
1:A:1272:G:H5'	14:N:34:VAL:HG11	1.91	0.52
22:a:1433:A:H2'	22:a:1434:A:C8	2.45	0.52
22:a:2700:A:H2'	22:a:2701:U:C6	2.45	0.52
44:w:43:GLU:CD	44:w:45:ARG:HE	2.17	0.52
1:A:236:A:H5''	17:Q:44:LEU:HD21	1.91	0.52
1:A:613:C:H2'	1:A:614:C:C6	2.45	0.52
1:A:1463:U:H2'	1:A:1464:U:C6	2.45	0.52
4:D:48:LEU:HD22	4:D:52:GLY:HA3	1.90	0.52
19:S:16:LEU:O	19:S:20:GLU:HG2	2.10	0.52
22:a:1568:G:N7	24:c:28:LYS:NZ	2.58	0.52
22:a:2795:C:H2'	22:a:2796:U:C6	2.45	0.52
1:A:76:G:N2	1:A:93:U:O2	2.30	0.52
22:a:1746:A:H2'	22:a:1747:U:C6	2.44	0.52
22:a:2273:A:H2'	22:a:2274:A:C8	2.45	0.52
26:e:61:ARG:HH12	26:e:64:GLY:HA3	1.75	0.52
29:h:34:GLY:O	29:h:36:ALA:N	2.42	0.52
1:A:736:C:H2'	1:A:737:C:C6	2.44	0.51
6:F:96:VAL:HG12	6:F:96:VAL:O	2.09	0.51
22:a:851:C:H2'	22:a:852:U:C6	2.45	0.51
22:a:2096:C:H2'	22:a:2097:A:C8	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:2455:G:H2'	22:a:2456:C:C6	2.45	0.51
33:l:64:TRP:HB2	33:l:104:GLU:HB2	1.91	0.51
10:J:80:THR:O	10:J:83:THR:OG1	2.20	0.51
1:A:736:C:H2'	1:A:737:C:H6	1.76	0.51
1:A:1412:C:H2'	1:A:1413:A:C8	2.45	0.51
22:a:172:A:H2'	22:a:173:A:H8	1.75	0.51
22:a:182:A:H2'	22:a:183:C:H6	1.76	0.51
22:a:742:A:H2'	22:a:743:A:H8	1.73	0.51
22:a:1248:G:OP1	26:e:44:ARG:NH1	2.42	0.51
22:a:1315:C:O2'	22:a:1392:A:N3	2.39	0.51
22:a:1405:U:H2'	22:a:1406:U:C6	2.46	0.51
22:a:2847:U:H2'	22:a:2848:G:O4'	2.09	0.51
18:R:9:LYS:NZ	18:R:43:ARG:O	2.43	0.51
22:a:1889:A:H2'	22:a:1890:A:C8	2.46	0.51
1:A:217:C:H2'	1:A:218:U:C6	2.46	0.51
2:B:154:MET:HE1	2:B:158:PRO:HG3	1.92	0.51
3:C:24:ALA:HB1	3:C:28:GLU:HG2	1.92	0.51
13:M:80:LEU:HD11	13:M:87:ARG:NH2	2.25	0.51
22:a:597:G:O2'	32:k:11:GLY:O	2.27	0.51
22:a:813:U:H2'	22:a:814:C:C6	2.45	0.51
22:a:2522:U:O2'	22:a:2647:U:OP1	2.20	0.51
37:p:49:ASP:HA	37:p:52:GLN:HB2	1.92	0.51
2:B:68:LEU:HD11	2:B:92:VAL:HG23	1.91	0.51
22:a:507:A:H5''	22:a:508:A:H3'	1.93	0.51
22:a:594:U:H2'	22:a:595:C:C6	2.46	0.51
22:a:2086:U:H2'	22:a:2087:G:C8	2.46	0.51
28:g:2:SER:OG	28:g:3:ARG:N	2.44	0.51
1:A:161:A:H2'	1:A:162:A:C8	2.45	0.51
1:A:613:C:OP1	4:D:81:ARG:NH1	2.39	0.51
1:A:932:C:H5''	7:G:4:ARG:NE	2.26	0.51
1:A:979:C:O2	14:N:59:ARG:NH1	2.44	0.51
6:F:18:VAL:HA	6:F:21:MET:HE2	1.91	0.51
22:a:125:A:OP2	49:1:19:ARG:NH1	2.35	0.51
22:a:1443:U:H2'	22:a:1444:G:H8	1.75	0.51
22:a:1538:G:H2'	22:a:1539:U:O4'	2.10	0.51
22:a:2700:A:H2'	22:a:2701:U:H6	1.76	0.51
31:j:38:ILE:HD11	31:j:112:PHE:CZ	2.45	0.51
40:s:20:ALA:O	40:s:24:MET:HG2	2.10	0.51
1:A:604:G:H2'	1:A:605:U:O4'	2.11	0.51
2:B:131:LYS:HE2	2:B:135:LEU:HD21	1.93	0.51
11:K:67:ALA:HB2	11:K:96:THR:HG23	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:75:ILE:HA	16:P:78:VAL:HG12	1.92	0.51
22:a:2848:G:O2'	22:a:2867:G:N2	2.36	0.51
47:z:9:THR:H	47:z:12:LYS:CE	2.24	0.51
13:M:16:VAL:O	13:M:20:THR:HG23	2.11	0.51
22:a:1664:A:N3	31:j:67:LYS:NZ	2.59	0.51
28:g:17:VAL:HG12	28:g:26:ILE:HG12	1.93	0.51
28:g:33:LEU:HD13	28:g:75:MET:HB2	1.93	0.51
53:V:54:5MU:H2'	53:V:55:PSU:C6	2.45	0.51
1:A:1328:C:H5''	13:M:28:THR:HG21	1.92	0.50
10:J:85:ASP:HB3	10:J:89:ARG:HH12	1.74	0.50
22:a:839:U:H2'	22:a:840:C:C6	2.46	0.50
22:a:2064:C:H2'	22:a:2065:C:C6	2.46	0.50
22:a:2845:U:H5''	36:o:52:ASN:O	2.10	0.50
45:x:10:SER:HB3	45:x:13:GLU:CD	2.36	0.50
55:X:23:A:H2'	55:X:24:G:C8	2.46	0.50
22:a:588:U:H2'	22:a:589:U:C6	2.46	0.50
22:a:640:C:H2'	22:a:641:U:H6	1.76	0.50
1:A:68:G:H22	1:A:101:A:H2	1.57	0.50
1:A:1465:A:H2'	1:A:1466:C:C6	2.47	0.50
22:a:78:U:H2'	22:a:79:C:C6	2.46	0.50
22:a:157:C:H2'	22:a:158:U:O4'	2.11	0.50
22:a:414:C:H2'	22:a:415:A:C8	2.47	0.50
22:a:414:C:H2'	22:a:415:A:H8	1.76	0.50
22:a:1563:U:H2'	22:a:1564:C:C6	2.47	0.50
1:A:1024:G:H2'	1:A:1025:U:O4'	2.11	0.50
3:C:47:LEU:HB3	3:C:50:ALA:HB3	1.93	0.50
7:G:5:ARG:NH2	7:G:7:ILE:HB	2.26	0.50
22:a:1386:C:H2'	22:a:1387:A:C8	2.47	0.50
1:A:461:A:H2'	1:A:462:G:C8	2.46	0.50
1:A:769:G:H4'	1:A:1513:A:H4'	1.92	0.50
2:B:23:TRP:HZ3	2:B:25:PRO:HA	1.77	0.50
3:C:121:THR:HA	3:C:124:LEU:HD12	1.94	0.50
22:a:1856:U:H2'	22:a:1857:G:O4'	2.11	0.50
22:a:2096:C:H2'	22:a:2097:A:H8	1.76	0.50
22:a:2567:G:H2'	22:a:2568:U:C6	2.47	0.50
46:y:39:GLU:OE1	46:y:41:THR:HG23	2.12	0.50
22:a:2395:C:H2'	22:a:2396:G:O4'	2.10	0.50
26:e:6:LYS:HD3	26:e:141:MET:HE3	1.94	0.50
13:M:34:LEU:HD23	13:M:56:LEU:HD11	1.93	0.50
22:a:150:U:H2'	22:a:151:C:C6	2.47	0.50
1:A:154:U:H2'	1:A:155:A:H8	1.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1071:C:H2'	1:A:1072:G:H8	1.77	0.50
14:N:50:THR:HG22	14:N:50:THR:O	2.11	0.50
1:A:451:A:H61	1:A:481:G:H5'	1.77	0.50
7:G:74:GLU:OE1	7:G:95:ARG:NH2	2.33	0.50
22:a:577:G:O2'	22:a:1254:A:OP1	2.30	0.50
1:A:177:G:OP2	1:A:177:G:N2	2.41	0.49
1:A:880:C:OP1	12:L:9:ARG:NH1	2.36	0.49
22:a:351:C:H2'	22:a:352:A:H8	1.77	0.49
22:a:582:A:H2'	22:a:583:G:C8	2.46	0.49
22:a:645:C:H2'	22:a:647:G:C8	2.47	0.49
22:a:703:U:H2'	22:a:704:G:O4'	2.12	0.49
25:d:6:GLY:O	25:d:201:LEU:N	2.40	0.49
1:A:662:U:H2'	1:A:663:A:C8	2.48	0.49
1:A:918:A:H2'	1:A:919:A:C8	2.47	0.49
1:A:996:A:H2'	1:A:997:U:C6	2.47	0.49
22:a:580:U:H2'	22:a:581:C:C6	2.47	0.49
22:a:1475:G:O2'	22:a:1514:G:O6	2.29	0.49
28:g:98:VAL:HG22	28:g:103:ILE:HG12	1.94	0.49
28:g:163:ARG:HG2	28:g:164:TYR:O	2.12	0.49
38:q:2:TYR:CD1	38:q:13:ARG:HD2	2.48	0.49
1:A:39:G:O6	1:A:547:A:H5''	2.13	0.49
1:A:1142:G:H2'	1:A:1143:G:O4'	2.12	0.49
1:A:1302:C:N4	13:M:14:HIS:HD2	2.09	0.49
7:G:16:PRO:HG2	7:G:44:TYR:OH	2.12	0.49
9:I:30:ILE:HG13	9:I:65:ILE:HB	1.94	0.49
16:P:36:VAL:HB	16:P:53:ASP:HB3	1.94	0.49
22:a:146:A:H2'	22:a:147:C:C6	2.47	0.49
22:a:920:A:OP1	46:y:19:LYS:NZ	2.45	0.49
22:a:1263:U:OP1	47:z:13:ARG:NH1	2.44	0.49
22:a:2478:A:H5'	51:3:32:LYS:HE2	1.93	0.49
23:b:66:A:OP2	23:b:108:A:N6	2.34	0.49
31:j:105:ARG:NH2	36:o:41:GLN:OE1	2.39	0.49
1:A:1314:C:H2'	1:A:1315:U:C6	2.47	0.49
4:D:95:GLU:HB3	4:D:186:PRO:HG2	1.94	0.49
13:M:24:GLY:O	13:M:29:ARG:NH1	2.45	0.49
22:a:849:A:H2'	22:a:850:U:C6	2.47	0.49
22:a:909:A:H2'	22:a:912:C:C5	2.48	0.49
22:a:1292:G:H2'	22:a:1293:C:C6	2.47	0.49
11:K:20:VAL:HG22	11:K:83:GLU:HB2	1.94	0.49
11:K:65:VAL:O	11:K:69:ARG:HG2	2.12	0.49
44:w:59:ILE:HD13	44:w:67:VAL:HG21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:4:58:ASP:OD1	52:4:59:ARG:N	2.46	0.49
1:A:626:G:O2'	16:P:51:ARG:NH2	2.43	0.49
1:A:1225:A:H2'	1:A:1226:C:C5	2.48	0.49
1:A:1410:A:H2'	1:A:1411:C:H6	1.77	0.49
8:H:59:LEU:HD12	8:H:60:GLU:H	1.78	0.49
22:a:5:A:H2'	22:a:6:A:C8	2.47	0.49
22:a:901:C:H2'	22:a:902:C:O4'	2.12	0.49
22:a:1412:U:H2'	22:a:1413:A:H8	1.77	0.49
30:i:95:ARG:HD2	30:i:96:ARG:HH11	1.78	0.49
55:X:72:C:H2'	55:X:73:A:O4'	2.12	0.49
1:A:155:A:H2'	1:A:156:C:H6	1.76	0.49
18:R:54:GLN:HA	18:R:57:ARG:NH1	2.27	0.49
19:S:12:ASP:OD2	19:S:35:SER:OG	2.18	0.49
19:S:30:PRO:HA	19:S:48:THR:HG23	1.95	0.49
22:a:419:U:H2'	22:a:420:C:C6	2.47	0.49
22:a:851:C:H2'	22:a:852:U:H6	1.78	0.49
30:i:96:ARG:O	30:i:96:ARG:HG2	2.12	0.49
1:A:677:U:O2	1:A:777:A:O2'	2.27	0.49
1:A:1162:C:H2'	1:A:1163:A:H8	1.78	0.49
22:a:668:A:H2'	22:a:670:A:H62	1.78	0.49
22:a:1526:C:H2'	22:a:1527:G:O4'	2.13	0.49
22:a:2723:C:H2'	22:a:2724:U:O4'	2.13	0.49
25:d:4:LEU:HD22	25:d:32:ASN:OD1	2.13	0.49
10:J:6:ILE:HG12	10:J:76:ILE:HG12	1.95	0.49
22:a:886:A:H2'	22:a:887:U:O4'	2.12	0.49
24:c:107:PRO:HD2	24:c:110:LEU:HD13	1.95	0.49
27:f:106:ILE:HG23	52:4:34:LEU:HD11	1.95	0.49
28:g:69:ARG:NH1	28:g:73:ASN:HB2	2.28	0.49
32:k:2:ARG:H	32:k:5:THR:CG2	2.26	0.49
1:A:1317:C:O2	19:S:37:ARG:NH2	2.46	0.49
3:C:86:LYS:HA	3:C:89:LYS:HE2	1.94	0.49
22:a:640:C:H2'	22:a:641:U:C6	2.47	0.49
22:a:1198:U:H2'	22:a:1199:U:C6	2.48	0.49
52:4:13:THR:OG1	52:4:23:LYS:NZ	2.46	0.49
55:X:11:C:H2'	55:X:12:U:C6	2.47	0.49
2:B:115:LYS:O	2:B:119:THR:HG23	2.13	0.48
22:a:956:G:O6	33:l:14:LYS:NZ	2.45	0.48
22:a:1197:G:H2'	22:a:1198:U:H6	1.78	0.48
22:a:1796:U:H2'	22:a:1797:G:C8	2.48	0.48
18:R:37:GLY:O	18:R:63:ARG:NH2	2.41	0.48
22:a:1432:G:H2'	22:a:1433:A:C8	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:s:30:ILE:HG13	40:s:85:VAL:HG13	1.95	0.48
22:a:543:G:H2'	22:a:544:C:C6	2.48	0.48
22:a:1682:G:H2'	22:a:1683:U:C6	2.47	0.48
22:a:2788:C:H2'	22:a:2789:C:C6	2.49	0.48
39:r:86:MET:HE3	39:r:96:ILE:HG13	1.96	0.48
55:X:23:A:H2'	55:X:24:G:H8	1.78	0.48
1:A:513:C:H2'	1:A:514:C:C6	2.48	0.48
1:A:784:A:H2'	1:A:785:G:C8	2.49	0.48
4:D:126:ASN:ND2	4:D:141:ASP:OD1	2.46	0.48
41:t:28:VAL:HG22	41:t:34:VAL:HG12	1.94	0.48
6:F:98:GLU:HG2	6:F:99:ALA:H	1.78	0.48
7:G:56:LYS:HG3	7:G:57:SER:H	1.77	0.48
14:N:48:LEU:HA	14:N:51:LEU:HD13	1.95	0.48
22:a:967:U:H2'	22:a:968:C:C6	2.49	0.48
23:b:106:G:H2'	23:b:107:G:O4'	2.12	0.48
36:o:91:ALA:HB2	36:o:113:ARG:HB3	1.95	0.48
5:E:11:LEU:HD12	5:E:68:ARG:HH12	1.78	0.48
9:I:28:ILE:HG12	9:I:63:LEU:HD12	1.96	0.48
11:K:83:GLU:OE1	11:K:109:ASN:HB2	2.12	0.48
22:a:24:G:O2'	39:r:78:GLU:O	2.32	0.48
22:a:2025:C:H2'	22:a:2026:U:C6	2.48	0.48
22:a:2799:A:O2'	22:a:2800:A:H5''	2.13	0.48
25:d:32:ASN:HD22	25:d:52:THR:HB	1.79	0.48
28:g:60:ASP:O	28:g:62:TRP:N	2.46	0.48
36:o:78:SER:OG	36:o:80:VAL:HG12	2.14	0.48
36:o:91:ALA:N	36:o:111:LYS:O	2.46	0.48
55:X:3:G:H1	55:X:70:C:H5	1.61	0.48
1:A:1530:G:H2'	1:A:1531:A:C8	2.49	0.48
2:B:83:ALA:HB2	2:B:214:LEU:HB3	1.94	0.48
3:C:19:ASN:O	3:C:40:ARG:NH2	2.47	0.48
11:K:16:VAL:HG12	11:K:18:ASP:H	1.78	0.48
22:a:927:A:H2'	22:a:928:A:C8	2.49	0.48
22:a:1179:G:H2'	22:a:1180:U:H6	1.78	0.48
22:a:1508:A:O2'	22:a:1509:A:O4'	2.27	0.48
22:a:2300:C:H2'	22:a:2301:C:H6	1.79	0.48
31:j:64:ARG:NH1	31:j:102:PRO:O	2.39	0.48
35:n:88:LYS:NZ	35:n:116:GLN:OE1	2.47	0.48
47:z:29:SER:OG	47:z:40:ARG:HD2	2.13	0.48
1:A:466:A:H2'	1:A:468:A:C8	2.49	0.48
1:A:1435:G:H2'	1:A:1436:U:C6	2.49	0.48
4:D:87:GLY:HA3	4:D:197:GLU:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:517:C:OP1	47:z:13:ARG:NH2	2.46	0.48
22:a:934:U:H2'	22:a:935:C:C6	2.48	0.48
22:a:1156:A:C8	37:p:51:ARG:HG2	2.49	0.48
22:a:2636:C:H2'	22:a:2637:U:C6	2.49	0.48
1:A:62:U:H2'	1:A:63:C:C6	2.49	0.48
1:A:91:U:H2'	1:A:92:U:C6	2.49	0.48
1:A:713:G:H2'	1:A:714:G:C8	2.48	0.48
5:E:34:THR:HG22	5:E:52:LYS:HG2	1.95	0.48
13:M:10:PRO:HG2	13:M:13:LYS:HE3	1.96	0.48
22:a:64:A:H2'	22:a:65:U:C6	2.48	0.48
22:a:581:C:H2'	22:a:582:A:C8	2.48	0.48
22:a:609:A:H2'	22:a:610:C:O4'	2.14	0.48
22:a:1199:U:H2'	22:a:1200:C:C6	2.49	0.48
22:a:2070:A:H2'	22:a:2071:A:C8	2.49	0.48
28:g:37:LEU:HD21	28:g:72:LEU:HD11	1.96	0.48
37:p:104:VAL:HG11	38:q:45:GLU:HG3	1.96	0.48
53:V:17:G:H1	53:V:55:PSU:HN3	1.61	0.48
1:A:335:C:H2'	1:A:336:A:H8	1.79	0.48
1:A:1442:G:H2'	1:A:1443:C:C6	2.48	0.48
10:J:25:ILE:HD11	10:J:92:LEU:HD11	1.95	0.48
22:a:861:A:H2'	22:a:862:G:O4'	2.14	0.48
22:a:2804:U:H2'	22:a:2805:C:H6	1.79	0.48
27:f:31:VAL:O	27:f:96:MET:HE1	2.14	0.48
1:A:136:C:O2'	16:P:64:GLY:O	2.21	0.47
1:A:323:U:H2'	1:A:324:G:O4'	2.13	0.47
1:A:744:C:H2'	1:A:745:G:H8	1.79	0.47
1:A:1442:G:H2'	1:A:1443:C:H6	1.79	0.47
7:G:28:ASN:OD1	7:G:36:LYS:HE2	2.14	0.47
16:P:52:LEU:HD12	16:P:57:ILE:HD11	1.94	0.47
22:a:476:G:H4'	22:a:502:A:N1	2.29	0.47
22:a:1187:G:H5''	38:q:83:TYR:CE1	2.49	0.47
22:a:1482:G:H1'	22:a:1509:A:N1	2.29	0.47
22:a:1809:A:H2'	22:a:1810:A:C8	2.49	0.47
1:A:33:A:H2'	1:A:34:C:C6	2.49	0.47
1:A:1010:U:H2'	1:A:1011:C:H6	1.79	0.47
6:F:38:ARG:HH22	18:R:24:LYS:HE3	1.79	0.47
10:J:59:LYS:HE2	10:J:62:ARG:NH2	2.29	0.47
13:M:87:ARG:HG2	13:M:97:VAL:HG13	1.95	0.47
17:Q:20:SER:HB3	17:Q:71:LYS:NZ	2.28	0.47
22:a:643:A:N7	22:a:644:A:C6	2.82	0.47
22:a:756:A:H2'	22:a:757:G:O4'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:1902:C:H4'	24:c:242:LYS:O	2.15	0.47
22:a:1939:5MU:OP1	22:a:2604:PSU:O2'	2.32	0.47
37:p:58:ARG:HA	37:p:61:TRP:CE3	2.49	0.47
44:w:55:GLY:O	44:w:59:ILE:HG12	2.14	0.47
1:A:543:U:OP1	4:D:14:ARG:NE	2.43	0.47
7:G:49:THR:O	7:G:52:GLN:HG3	2.13	0.47
16:P:52:LEU:HD21	16:P:74:LEU:HD13	1.97	0.47
20:T:27:MET:O	20:T:30:THR:HG22	2.14	0.47
22:a:95:A:H4'	45:x:38:GLN:O	2.14	0.47
22:a:288:U:H2'	22:a:289:G:C8	2.49	0.47
22:a:1179:G:H2'	22:a:1180:U:C6	2.48	0.47
22:a:1469:A:H2'	22:a:1470:A:H8	1.79	0.47
22:a:1704:C:H2'	22:a:1705:A:C8	2.50	0.47
33:l:136:MET:O	42:u:79:ARG:NH2	2.33	0.47
2:B:14:VAL:HG23	2:B:213:TYR:OH	2.14	0.47
18:R:21:ILE:HG21	18:R:55:LEU:HA	1.96	0.47
19:S:50:ALA:HB1	19:S:57:HIS:HB3	1.97	0.47
22:a:235:U:H2'	22:a:236:C:H6	1.79	0.47
40:s:15:HIS:HB3	40:s:31:VAL:HG12	1.96	0.47
1:A:763:G:H2'	1:A:764:C:C6	2.49	0.47
5:E:162:GLU:H	5:E:162:GLU:CD	2.22	0.47
8:H:87:LYS:HD2	8:H:91:GLU:HG3	1.97	0.47
18:R:26:ILE:HD13	18:R:29:LEU:HD12	1.95	0.47
22:a:1539:U:H2'	22:a:1540:G:C8	2.49	0.47
22:a:1796:U:H2'	22:a:1797:G:H8	1.79	0.47
22:a:2392:A:O2'	32:k:60:ARG:O	2.32	0.47
28:g:35:ARG:HB3	28:g:75:MET:HE1	1.96	0.47
31:j:40:LYS:HD3	31:j:58:LEU:O	2.14	0.47
1:A:107:G:N1	20:T:6:SER:OG	2.36	0.47
1:A:467:U:H3'	1:A:468:A:C5'	2.44	0.47
1:A:579:A:H2'	1:A:580:C:C6	2.50	0.47
1:A:592:G:H2'	1:A:593:U:H6	1.79	0.47
1:A:1266:G:N2	1:A:1269:A:OP2	2.38	0.47
2:B:89:GLN:OE1	2:B:89:GLN:HA	2.15	0.47
4:D:25:VAL:HB	4:D:161:LEU:HD21	1.96	0.47
6:F:38:ARG:HH12	6:F:100:SER:HB3	1.78	0.47
10:J:15:HIS:HB3	10:J:70:HIS:CE1	2.50	0.47
22:a:760:G:H2'	22:a:761:A:O4'	2.15	0.47
22:a:848:C:H2'	22:a:849:A:C8	2.47	0.47
22:a:898:C:H2'	22:a:899:A:O4'	2.15	0.47
22:a:1013:C:H2'	22:a:1014:A:C8	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:1443:U:H2'	22:a:1444:G:C8	2.49	0.47
22:a:2430:A:N3	22:a:2430:A:H2'	2.28	0.47
36:o:3:ASN:HA	36:o:6:LYS:HG2	1.97	0.47
1:A:384:G:H2'	1:A:385:C:H6	1.80	0.47
5:E:107:ALA:HB2	5:E:125:ALA:HB3	1.97	0.47
6:F:18:VAL:HG21	6:F:58:HIS:ND1	2.29	0.47
22:a:1417:C:O2'	22:a:1587:G:O2'	2.27	0.47
22:a:2246:G:H2'	22:a:2247:A:C8	2.49	0.47
1:A:41:G:H2'	1:A:42:G:H8	1.79	0.47
1:A:977:A:OP1	14:N:61:ARG:NH2	2.42	0.47
10:J:9:ARG:HB2	10:J:99:GLN:HB2	1.97	0.47
19:S:49:ILE:O	19:S:60:VAL:HG12	2.14	0.47
22:a:355:U:H2'	22:a:356:G:H8	1.78	0.47
22:a:1728:C:H2'	22:a:1729:U:H5''	1.96	0.47
22:a:2625:G:H2'	22:a:2626:C:O4'	2.15	0.47
25:d:121:THR:CG2	25:d:161:MET:HA	2.43	0.47
52:4:8:LYS:HG2	52:4:10:GLU:OE1	2.15	0.47
22:a:348:A:H2'	22:a:349:U:C6	2.50	0.47
22:a:357:C:H2'	22:a:358:U:C6	2.50	0.47
22:a:2806:C:H2'	22:a:2807:U:O4'	2.15	0.47
26:e:170:ARG:HH12	26:e:179:SER:CB	2.27	0.47
1:A:72:A:H2'	1:A:72:A:N3	2.29	0.46
1:A:1391:U:H2'	1:A:1392:G:H8	1.79	0.46
11:K:89:PRO:HB3	21:U:32:VAL:HG21	1.96	0.46
20:T:44:LYS:HD2	20:T:87:ALA:HA	1.98	0.46
22:a:2014:A:H2'	22:a:2015:A:C8	2.50	0.46
22:a:2097:A:H2'	22:a:2098:U:C6	2.50	0.46
23:b:112:G:H2'	23:b:113:C:C6	2.50	0.46
24:c:51:THR:HG23	24:c:54:ILE:HD12	1.96	0.46
1:A:1031:C:O2'	1:A:1032:G:N2	2.49	0.46
1:A:1119:C:H2'	1:A:1120:C:H6	1.80	0.46
1:A:1404:C:H2'	1:A:1405:G:C8	2.50	0.46
13:M:55:THR:O	13:M:59:GLU:HG2	2.15	0.46
13:M:73:ILE:O	13:M:77:ILE:HG12	2.15	0.46
22:a:277:G:H4'	22:a:278:A:C4	2.51	0.46
22:a:302:C:H2'	22:a:303:G:H8	1.81	0.46
22:a:1509:A:HO2'	22:a:1510:G:H8	1.62	0.46
22:a:1704:C:H2'	22:a:1705:A:H8	1.80	0.46
22:a:2051:A:H8	22:a:2051:A:OP2	1.98	0.46
22:a:2327:A:H2'	22:a:2328:A:C8	2.50	0.46
1:A:590:U:H2'	1:A:591:U:C6	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:674:G:H2'	1:A:675:A:C8	2.50	0.46
1:A:950:U:H2'	1:A:951:G:C8	2.50	0.46
2:B:100:MET:O	2:B:104:TRP:HE3	1.98	0.46
21:U:11:PRO:O	21:U:14:VAL:HG22	2.16	0.46
22:a:1178:C:H2'	22:a:1179:G:C8	2.49	0.46
22:a:2020:A:H5'	47:z:9:THR:CG2	2.46	0.46
27:f:100:PHE:O	27:f:104:ILE:HG22	2.16	0.46
1:A:591:U:H2'	1:A:592:G:H8	1.80	0.46
1:A:1169:A:H2'	1:A:1170:A:C8	2.50	0.46
4:D:62:ARG:O	4:D:66:GLY:N	2.48	0.46
22:a:127:A:H62	49:1:46:LYS:HD2	1.80	0.46
22:a:1932:A:H2'	22:a:1933:G:O4'	2.15	0.46
24:c:231:PRO:O	24:c:242:LYS:HE3	2.16	0.46
32:k:62:PRO:HG2	50:2:25:LYS:HB3	1.97	0.46
1:A:131:A:H2'	1:A:132:C:H6	1.81	0.46
1:A:236:A:H2'	1:A:237:G:H8	1.81	0.46
1:A:529:G:H5'	1:A:530:G:OP2	2.15	0.46
26:e:168:ASP:OD2	26:e:170:ARG:NH2	2.49	0.46
42:u:77:VAL:HG23	42:u:89:ILE:HG12	1.97	0.46
1:A:983:A:H5'	1:A:984:C:OP2	2.15	0.46
1:A:1239:A:O2'	7:G:114:LYS:O	2.31	0.46
16:P:6:LEU:HB3	16:P:17:TYR:HB3	1.97	0.46
22:a:131:A:H2'	22:a:132:G:C8	2.51	0.46
55:X:75:C:H2'	55:X:76:A:C8	2.50	0.46
1:A:41:G:H2'	1:A:42:G:C8	2.50	0.46
1:A:80:A:H2'	1:A:81:A:H8	1.80	0.46
1:A:821:G:H2'	1:A:822:U:C6	2.51	0.46
22:a:721:A:H2'	22:a:722:A:C8	2.51	0.46
22:a:784:G:H5'	22:a:785:G:OP1	2.16	0.46
22:a:1794:A:H2'	22:a:1795:C:H6	1.78	0.46
30:i:69:ARG:O	30:i:90:GLU:HG3	2.16	0.46
5:E:95:PHE:CE2	5:E:97:GLN:HG3	2.51	0.46
14:N:9:ARG:HB3	14:N:13:ARG:NH1	2.31	0.46
22:a:235:U:H2'	22:a:236:C:C6	2.51	0.46
22:a:1496:A:H2'	22:a:1498:C:C5	2.51	0.46
22:a:2038:G:H2'	22:a:2039:U:O4'	2.15	0.46
31:j:66:LYS:HE2	31:j:66:LYS:HB3	1.65	0.46
40:s:64:LYS:HD2	40:s:64:LYS:N	2.30	0.46
52:4:3:LYS:HA	52:4:3:LYS:HD3	1.72	0.46
1:A:176:C:H2'	1:A:177:G:N3	2.31	0.46
1:A:695:A:H2'	1:A:696:A:C8	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1163:A:H2'	1:A:1164:G:C8	2.51	0.46
9:I:41:ARG:NH2	9:I:43:THR:OG1	2.48	0.46
14:N:38:ASP:OD1	14:N:39:GLU:N	2.49	0.46
22:a:2636:C:H2'	22:a:2637:U:H6	1.81	0.46
26:e:21:ARG:NH2	26:e:106:LYS:HD3	2.28	0.46
30:i:110:PRO:O	30:i:115:GLY:HA3	2.16	0.46
41:t:34:VAL:HG13	41:t:67:VAL:HG12	1.97	0.46
55:X:18:G:O2'	55:X:19:G:OP2	2.30	0.46
1:A:324:G:N1	1:A:327:A:OP2	2.48	0.46
1:A:1476:A:H2'	1:A:1477:U:C6	2.51	0.46
1:A:1513:A:H2'	1:A:1514:G:C8	2.51	0.46
6:F:11:HIS:ND1	6:F:12:PRO:HD2	2.31	0.46
8:H:39:VAL:HG21	8:H:110:VAL:HG12	1.97	0.46
13:M:81:MET:O	13:M:92:ARG:NH2	2.36	0.46
21:U:62:ARG:O	21:U:66:ARG:HG3	2.16	0.46
22:a:223:A:N1	22:a:407:G:O2'	2.45	0.46
22:a:372:G:O2'	44:w:54:LYS:HE3	2.16	0.46
22:a:589:U:H2'	22:a:590:A:C8	2.51	0.46
22:a:918:A:H5''	23:b:97:C:O2'	2.16	0.46
22:a:2419:U:OP2	50:2:33:LEU:HD13	2.16	0.46
22:a:2469:A:N6	22:a:2481:G:O2'	2.49	0.46
2:B:186:ILE:HD13	2:B:200:ILE:HB	1.98	0.45
5:E:11:LEU:HD12	5:E:68:ARG:NH1	2.31	0.45
8:H:34:VAL:HG13	8:H:49:PHE:CE2	2.50	0.45
22:a:550:C:H2'	22:a:551:G:H8	1.81	0.45
22:a:1656:C:H2'	22:a:1657:U:H6	1.81	0.45
39:r:23:LEU:HD22	47:z:24:ALA:HB2	1.96	0.45
1:A:539:A:H2'	1:A:540:G:C8	2.51	0.45
1:A:592:G:H2'	1:A:593:U:C6	2.50	0.45
1:A:1033:G:OP2	1:A:1033:G:N2	2.39	0.45
11:K:60:PRO:HD3	11:K:91:PRO:HB2	1.99	0.45
11:K:64:GLN:HG3	11:K:99:ALA:HB2	1.97	0.45
22:a:322:A:OP2	26:e:163:ASN:HB2	2.16	0.45
22:a:2901:C:N4	22:a:2902:C:H41	2.14	0.45
25:d:37:VAL:HG23	25:d:92:VAL:HG23	1.98	0.45
1:A:696:A:H2'	1:A:697:U:H6	1.81	0.45
2:B:162:PHE:CZ	2:B:217:VAL:HG11	2.52	0.45
22:a:2251:OMG:HM23	22:a:2251:OMG:H1'	1.72	0.45
27:f:36:LEU:HB2	27:f:89:VAL:HG12	1.98	0.45
30:i:96:ARG:HE	30:i:99:ARG:HG3	1.81	0.45
10:J:17:LEU:HD21	10:J:94:ALA:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:308:G:H2'	22:a:309:A:C8	2.51	0.45
22:a:2646:C:H2'	22:a:2647:U:O4'	2.17	0.45
27:f:175:PHE:O	27:f:177:PHE:N	2.49	0.45
37:p:107:THR:HA	37:p:110:VAL:HG12	1.98	0.45
50:2:32:ILE:HD12	50:2:35:LYS:HE2	1.98	0.45
22:a:166:U:H2'	22:a:167:A:H8	1.80	0.45
22:a:479:A:N3	22:a:481:G:H5''	2.32	0.45
40:s:64:LYS:HA	40:s:79:ASP:OD1	2.17	0.45
22:a:458:G:O2'	22:a:469:G:O6	2.34	0.45
22:a:2585:U:HO2'	22:a:2586:U:P	2.39	0.45
25:d:38:LYS:HE3	25:d:81:GLU:OE2	2.17	0.45
27:f:46:ASP:OD2	27:f:48:LYS:HB2	2.15	0.45
1:A:119:A:H4'	1:A:120:A:C4	2.52	0.45
1:A:440:C:C2	1:A:441:A:C8	3.04	0.45
14:N:88:ALA:HB2	14:N:96:LEU:HD23	1.99	0.45
22:a:729:G:C6	24:c:207:LYS:HB2	2.52	0.45
22:a:833:A:H2'	22:a:834:G:C8	2.52	0.45
22:a:1441:G:H2'	22:a:1442:U:C6	2.51	0.45
22:a:2047:C:H2'	22:a:2048:G:H8	1.82	0.45
24:c:76:ALA:HB2	24:c:96:TYR:CD2	2.52	0.45
45:x:12:GLU:HG2	45:x:13:GLU:N	2.31	0.45
55:X:8:U:O2'	55:X:21:A:N1	2.42	0.45
1:A:993:G:O2'	1:A:994:A:N7	2.50	0.45
9:I:12:ARG:HG3	9:I:13:LYS:H	1.81	0.45
10:J:5:ARG:HH12	10:J:7:ARG:NE	2.15	0.45
22:a:247:G:H4'	22:a:386:G:C5	2.51	0.45
22:a:1199:U:H2'	22:a:1200:C:H6	1.82	0.45
22:a:1387:A:H2'	22:a:1388:G:H8	1.82	0.45
22:a:2557:G:H2'	22:a:2558:C:C6	2.52	0.45
24:c:5:LYS:HD2	24:c:17:VAL:HG22	1.98	0.45
22:a:17:G:H2'	22:a:18:U:C6	2.52	0.45
22:a:132:G:H2'	22:a:133:U:C6	2.51	0.45
22:a:624:C:O2'	22:a:657:U:OP1	2.33	0.45
22:a:832:U:H2'	22:a:833:A:H8	1.82	0.45
22:a:839:U:H2'	22:a:840:C:H6	1.82	0.45
22:a:1405:U:H2'	22:a:1406:U:H6	1.80	0.45
40:s:54:GLU:H	40:s:91:GLN:HG2	1.82	0.45
55:X:6:C:H4'	55:X:7:U:OP1	2.17	0.45
1:A:215:C:H2'	1:A:216:U:C6	2.52	0.45
1:A:381:C:H2'	1:A:382:A:O4'	2.17	0.45
1:A:1120:C:H2'	1:A:1121:U:H6	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:19:LEU:HD22	4:D:64:ILE:HG13	1.99	0.45
7:G:15:ASP:HB3	7:G:20:SER:H	1.82	0.45
7:G:71:PRO:O	7:G:96:ARG:HG2	2.17	0.45
8:H:48:ASP:OD1	8:H:49:PHE:N	2.50	0.45
9:I:88:MET:HE1	9:I:95:ARG:HA	1.98	0.45
14:N:42:TRP:O	14:N:45:VAL:HG12	2.17	0.45
22:a:150:U:H2'	22:a:151:C:H6	1.81	0.45
22:a:590:A:H2'	22:a:591:U:C6	2.52	0.45
22:a:1545:A:H2'	22:a:1546:G:O4'	2.17	0.45
22:a:1744:A:H3'	22:a:1745:A:H8	1.82	0.45
22:a:2615:U:C2	47:z:4:GLN:HA	2.51	0.45
22:a:2898:U:O2	30:i:134:ALA:HB1	2.16	0.45
24:c:145:GLU:HB2	24:c:188:CYS:HB3	1.99	0.45
27:f:114:PHE:HZ	27:f:176:PRO:HB2	1.82	0.45
32:k:108:ALA:HB3	32:k:125:LEU:HD22	1.99	0.45
2:B:105:LYS:HA	2:B:108:ARG:HD3	1.98	0.44
22:a:1378:A:O2'	22:a:1380:G:N7	2.50	0.44
22:a:1387:A:H2'	22:a:1388:G:C8	2.52	0.44
40:s:59:ASN:O	40:s:84:TYR:N	2.49	0.44
1:A:580:C:H2'	1:A:581:G:O4'	2.18	0.44
1:A:1123:U:O2'	1:A:1124:G:H5'	2.16	0.44
1:A:1248:A:H2'	1:A:1249:C:C6	2.52	0.44
1:A:1530:G:N7	21:U:46:LYS:HE3	2.31	0.44
6:F:2:ARG:HD3	6:F:91:ARG:CZ	2.47	0.44
17:Q:47:HIS:NE2	17:Q:49:GLU:HG2	2.33	0.44
20:T:42:GLY:HA2	20:T:86:LEU:HD11	1.98	0.44
22:a:739:A:H1'	22:a:740:C:H5	1.83	0.44
22:a:1242:U:H2'	22:a:1243:C:C6	2.52	0.44
36:o:51:ARG:CZ	36:o:53:ARG:HG3	2.47	0.44
1:A:171:A:H2'	1:A:172:A:C8	2.53	0.44
1:A:1315:U:H2'	1:A:1316:G:O4'	2.17	0.44
2:B:188:ASP:CG	2:B:189:THR:N	2.70	0.44
6:F:74:LEU:O	6:F:77:THR:OG1	2.28	0.44
20:T:29:ARG:O	20:T:33:LYS:HG2	2.17	0.44
22:a:359:G:H2'	22:a:360:U:C6	2.52	0.44
22:a:1730:C:O2'	22:a:1731:G:O5'	2.35	0.44
22:a:1946:U:H2'	22:a:1947:C:C6	2.52	0.44
22:a:2305:U:H5''	27:f:131:GLY:HA3	1.99	0.44
22:a:2849:U:H4'	22:a:2868:A:C2	2.52	0.44
1:A:744:C:H2'	1:A:745:G:C8	2.52	0.44
1:A:865:A:H2'	1:A:866:C:C6	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:97:LEU:HD11	2:B:147:SER:OG	2.17	0.44
6:F:42:TRP:NE1	6:F:102:MET:HB2	2.32	0.44
7:G:79:ARG:NH2	7:G:82:GLY:HA2	2.31	0.44
7:G:113:ASP:O	7:G:119:ARG:NH2	2.43	0.44
18:R:21:ILE:HG13	18:R:54:GLN:HB2	1.99	0.44
21:U:39:GLU:OE1	21:U:47:ARG:NH1	2.51	0.44
22:a:1490:A:O2'	22:a:1491:G:H5'	2.17	0.44
22:a:1593:A:H2'	22:a:1594:U:C6	2.52	0.44
28:g:28:GLY:HA3	28:g:79:VAL:HG22	2.00	0.44
1:A:575:G:O2'	1:A:821:G:OP2	2.33	0.44
1:A:663:A:H5'	1:A:836:G:OP1	2.18	0.44
1:A:1162:C:H2'	1:A:1163:A:C8	2.53	0.44
17:Q:8:LEU:HD23	17:Q:8:LEU:HA	1.83	0.44
22:a:813:U:H2'	22:a:814:C:H6	1.81	0.44
22:a:1727:C:H2'	22:a:1728:C:C6	2.52	0.44
26:e:106:LYS:HG3	26:e:200:LEU:HD13	1.99	0.44
32:k:77:ILE:CD1	32:k:108:ALA:HB1	2.48	0.44
55:X:26:M2G:H3'	55:X:27:A:H8	1.81	0.44
11:K:85:MET:HE3	11:K:113:VAL:HG11	2.00	0.44
22:a:184:C:H2'	22:a:185:G:H8	1.82	0.44
22:a:518:G:H2'	22:a:519:U:C6	2.52	0.44
22:a:581:C:H2'	22:a:582:A:H8	1.82	0.44
22:a:930:G:H1'	46:y:25:LEU:HD21	1.99	0.44
23:b:15:A:H3'	23:b:16:G:H8	1.83	0.44
24:c:174:LEU:HD11	24:c:184:VAL:HG11	1.99	0.44
27:f:58:ALA:HB2	27:f:65:PRO:HD3	1.99	0.44
36:o:100:LEU:HD11	36:o:110:ILE:HD11	2.00	0.44
37:p:107:THR:O	37:p:111:GLU:HG2	2.18	0.44
46:y:41:THR:HB	46:y:42:PRO:HD2	2.00	0.44
6:F:73:GLU:O	6:F:77:THR:HG23	2.18	0.44
22:a:3:U:H2'	22:a:4:U:C6	2.52	0.44
22:a:638:G:H2'	22:a:639:U:C6	2.53	0.44
22:a:994:C:OP2	37:p:50:ARG:HG2	2.17	0.44
22:a:1677:A:H2'	22:a:1678:A:C8	2.53	0.44
22:a:1736:U:H2'	22:a:1737:G:O4'	2.18	0.44
28:g:78:GLY:HA2	28:g:82:GLY:HA2	2.00	0.44
36:o:18:PRO:HG3	36:o:84:ILE:O	2.17	0.44
1:A:608:A:H2'	1:A:609:A:O4'	2.18	0.44
2:B:162:PHE:CE1	2:B:217:VAL:HG11	2.53	0.44
3:C:179:ARG:HG2	3:C:207:ILE:HA	2.00	0.44
22:a:391:A:H1'	22:a:411:G:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:500:G:N1	22:a:503:A:OP2	2.47	0.44
22:a:594:U:H2'	22:a:595:C:H6	1.82	0.44
22:a:743:A:O2'	22:a:1659:G:OP1	2.29	0.44
22:a:849:A:H2'	22:a:850:U:H6	1.81	0.44
22:a:1649:G:O2'	34:m:106:ASP:OD2	2.19	0.44
22:a:1916:A:H2'	22:a:1917:PSU:O4'	2.18	0.44
26:e:111:GLU:OE1	26:e:114:ARG:NH2	2.38	0.44
28:g:72:LEU:HA	28:g:75:MET:HG2	2.00	0.44
31:j:20:MET:SD	31:j:44:LYS:HE3	2.58	0.44
39:r:43:ALA:HA	39:r:46:LEU:HD12	1.98	0.44
1:A:16:A:N1	1:A:919:A:H2	2.15	0.44
1:A:1507:A:H2'	1:A:1508:A:C8	2.53	0.44
22:a:231:A:H2'	22:a:232:G:O4'	2.18	0.44
22:a:244:A:H2'	22:a:245:G:O4'	2.18	0.44
22:a:645:C:H2'	22:a:647:G:N7	2.33	0.44
22:a:1028:A:N6	22:a:1125:G:H2'	2.33	0.44
22:a:2420:C:H4'	48:0:54:ILE:HG21	2.00	0.44
22:a:2752:C:H2'	22:a:2753:A:O4'	2.18	0.44
47:z:9:THR:H	47:z:12:LYS:HE3	1.82	0.44
53:V:9:C:O2'	53:V:10:G:N7	2.45	0.44
1:A:219:U:H2'	1:A:220:G:H8	1.83	0.43
1:A:232:G:H1'	1:A:262:A:N1	2.33	0.43
1:A:349:A:O2'	1:A:350:G:H5'	2.18	0.43
1:A:519:C:HO2'	1:A:520:A:P	2.41	0.43
1:A:868:C:H2'	1:A:869:G:O4'	2.18	0.43
2:B:76:ALA:HB1	2:B:210:VAL:HG11	2.00	0.43
2:B:87:CYS:HB3	2:B:218:ALA:HB1	2.00	0.43
4:D:139:PRO:HD3	4:D:185:LYS:HE2	2.00	0.43
22:a:5:A:H2'	22:a:6:A:H8	1.83	0.43
22:a:538:A:H2'	22:a:539:G:O4'	2.18	0.43
22:a:580:U:H2'	22:a:581:C:H6	1.82	0.43
22:a:1266:G:O2'	22:a:2012:G:O6	2.31	0.43
33:l:106:ASP:OD1	33:l:107:GLY:N	2.47	0.43
34:m:103:ARG:HD2	34:m:106:ASP:OD1	2.18	0.43
36:o:6:LYS:O	36:o:10:GLN:HG2	2.18	0.43
40:s:53:VAL:HG11	40:s:93:LEU:HB2	2.00	0.43
42:u:2:PHE:CD1	42:u:50:MET:HE3	2.53	0.43
42:u:30:ILE:HG12	42:u:91:PHE:HB2	2.00	0.43
1:A:216:U:H2'	1:A:217:C:H6	1.83	0.43
1:A:591:U:OP1	8:H:31:LYS:HE3	2.17	0.43
1:A:831:A:H4'	2:B:21:ARG:NH2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1175:G:H2'	1:A:1176:A:C8	2.51	0.43
7:G:56:LYS:HB2	7:G:60:GLU:OE2	2.18	0.43
21:U:50:ALA:HA	21:U:53:VAL:HG12	1.99	0.43
22:a:709:U:H2'	22:a:710:U:C6	2.53	0.43
22:a:1572:A:H2'	22:a:1573:G:H8	1.83	0.43
22:a:2216:G:H2'	22:a:2217:G:H8	1.84	0.43
22:a:2698:U:H2'	22:a:2699:C:C6	2.53	0.43
24:c:72:ASP:OD2	24:c:119:GLY:HA2	2.18	0.43
1:A:192:A:H2'	1:A:193:C:C6	2.53	0.43
1:A:471:U:H2'	1:A:472:U:H6	1.84	0.43
1:A:553:A:H5''	12:L:21:VAL:HG21	2.00	0.43
3:C:153:VAL:HG12	3:C:198:VAL:HG22	2.00	0.43
10:J:11:LYS:HB2	10:J:97:ASP:OD1	2.18	0.43
13:M:19:LEU:HD23	13:M:19:LEU:HA	1.88	0.43
13:M:89:LEU:HB3	13:M:93:ARG:NH1	2.33	0.43
22:a:68:G:H2'	22:a:69:C:O4'	2.18	0.43
22:a:78:U:H2'	22:a:79:C:H6	1.82	0.43
22:a:1641:A:H2'	22:a:1642:G:O4'	2.19	0.43
22:a:1727:C:H2'	22:a:1728:C:H6	1.84	0.43
22:a:1790:C:H2'	22:a:1791:A:C5	2.52	0.43
40:s:4:GLU:OE1	40:s:49:LYS:NZ	2.36	0.43
1:A:34:C:H2'	1:A:35:G:H8	1.83	0.43
1:A:246:A:C2	1:A:282:A:C5	3.06	0.43
1:A:986:U:H2'	1:A:987:G:C8	2.53	0.43
1:A:1314:C:H2'	1:A:1315:U:H6	1.83	0.43
2:B:134:ALA:O	2:B:138:THR:HG23	2.19	0.43
2:B:213:TYR:O	2:B:217:VAL:HG22	2.19	0.43
9:I:118:LEU:HD22	9:I:124:ARG:HG2	1.99	0.43
13:M:92:ARG:HG2	22:a:888:C:C6	2.53	0.43
18:R:11:CYS:SG	18:R:48:ARG:HG3	2.58	0.43
22:a:2680:U:H5'	25:d:194:PRO:HA	2.00	0.43
22:a:2790:U:H5'	22:a:2893:A:H62	1.84	0.43
35:n:25:ARG:HG2	35:n:27:VAL:HG22	1.99	0.43
38:q:1:MET:HE3	38:q:101:ILE:HB	1.99	0.43
1:A:502:A:OP1	12:L:115:SER:OG	2.28	0.43
1:A:546:A:OP2	4:D:68:LEU:HD12	2.19	0.43
7:G:111:ARG:HB3	7:G:119:ARG:HG2	2.00	0.43
22:a:753:A:H2'	22:a:754:U:H6	1.84	0.43
22:a:1429:G:H2'	22:a:1430:G:H8	1.83	0.43
22:a:1710:G:H4'	22:a:2858:C:O2	2.18	0.43
25:d:121:THR:HB	25:d:127:PHE:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:u:47:VAL:O	42:u:51:GLN:HG3	2.19	0.43
1:A:74:A:H2'	1:A:75:G:H8	1.83	0.43
1:A:526:C:OP2	12:L:88:LYS:HE3	2.19	0.43
1:A:718:A:H5'	11:K:119:IAS:C	2.49	0.43
1:A:746:A:H2'	1:A:747:A:C8	2.53	0.43
1:A:855:U:OP2	1:A:871:U:N3	2.44	0.43
1:A:1312:G:H5'	19:S:5:LEU:HD11	1.99	0.43
3:C:150:LYS:HE3	3:C:201:TRP:CE3	2.53	0.43
5:E:85:VAL:HG11	5:E:147:MET:HG3	2.00	0.43
10:J:10:LEU:HB2	10:J:72:ARG:HB2	1.99	0.43
22:a:828:U:H2'	22:a:829:A:C8	2.54	0.43
22:a:852:U:H2'	22:a:853:C:C6	2.54	0.43
22:a:1800:C:H5'	24:c:146:MET:HE1	2.00	0.43
23:b:14:U:OP2	23:b:70:C:O2'	2.34	0.43
34:m:56:LYS:HE2	34:m:87:PHE:O	2.18	0.43
37:p:106:PHE:HA	37:p:109:LEU:HD12	2.01	0.43
1:A:160:A:H2'	1:A:161:A:O4'	2.18	0.43
1:A:182:A:N1	1:A:223:A:O2'	2.52	0.43
1:A:523:A:C2	12:L:88:LYS:HB3	2.53	0.43
16:P:19:VAL:HG13	16:P:36:VAL:HG23	2.01	0.43
22:a:12:U:O2	22:a:2626:C:H4'	2.19	0.43
22:a:1316:U:H2'	22:a:1317:G:C8	2.54	0.43
22:a:2246:G:H2'	22:a:2247:A:H8	1.84	0.43
33:l:36:VAL:HG12	42:u:82:TYR:HB2	2.00	0.43
36:o:99:TYR:O	36:o:102:GLU:HG2	2.18	0.43
41:t:72:ILE:HD13	41:t:103:ILE:HD13	2.01	0.43
55:X:67:G:H2'	55:X:68:G:C8	2.54	0.43
1:A:1236:A:H2'	1:A:1237:C:C6	2.53	0.43
9:I:7:TYR:OH	9:I:9:THR:HB	2.19	0.43
22:a:851:C:O2'	46:y:43:ALA:O	2.35	0.43
22:a:1278:C:H2'	22:a:1279:G:H8	1.83	0.43
22:a:1680:U:H2'	22:a:1681:G:O4'	2.19	0.43
22:a:1786:A:H1'	22:a:1938:A:N6	2.33	0.43
22:a:2530:A:N7	28:g:172:LYS:NZ	2.63	0.43
28:g:140:VAL:O	28:g:144:VAL:HG23	2.18	0.43
34:m:71:ARG:HD2	34:m:71:ARG:HA	1.79	0.43
1:A:1238:A:H2	1:A:1241:G:N3	2.17	0.43
1:A:1356:G:H2'	1:A:1357:A:H8	1.78	0.43
4:D:109:ALA:H	4:D:113:GLU:HG2	1.83	0.43
10:J:90:LEU:HB3	10:J:92:LEU:HD22	2.01	0.43
21:U:56:HIS:O	21:U:60:LEU:HD23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:278:A:C6	22:a:362:A:C8	3.07	0.43
28:g:72:LEU:HG	28:g:75:MET:HE2	2.01	0.43
49:1:42:LEU:HD23	49:1:42:LEU:HA	1.85	0.43
1:A:269:C:H2'	1:A:270:A:H8	1.84	0.43
1:A:312:C:H2'	1:A:313:A:C8	2.54	0.43
1:A:503:C:O2'	1:A:510:A:N1	2.46	0.43
1:A:625:U:H2'	1:A:626:G:C8	2.54	0.43
1:A:1233:G:H4'	9:I:119:ARG:HH12	1.84	0.43
1:A:1438:G:H5''	20:T:33:LYS:NZ	2.33	0.43
10:J:8:ILE:HA	10:J:99:GLN:O	2.19	0.43
22:a:103:A:H2'	22:a:104:A:O4'	2.18	0.43
22:a:208:C:H2'	22:a:209:C:C6	2.54	0.43
22:a:263:G:H2'	22:a:264:C:O4'	2.19	0.43
22:a:550:C:H2'	22:a:551:G:C8	2.54	0.43
22:a:1778:U:H2'	22:a:1784:A:N6	2.34	0.43
22:a:1799:G:O2'	24:c:180:GLU:OE2	2.33	0.43
25:d:32:ASN:ND2	25:d:52:THR:HB	2.33	0.43
25:d:152:PRO:HG3	25:d:156:PHE:CZ	2.54	0.43
50:2:32:ILE:O	50:2:32:ILE:HG13	2.19	0.43
1:A:35:G:H2'	1:A:36:C:C6	2.54	0.42
9:I:55:VAL:HG23	9:I:57:MET:HG3	2.01	0.42
21:U:4:ILE:HG13	21:U:19:PHE:HA	1.99	0.42
22:a:151:C:H2'	22:a:152:A:H8	1.83	0.42
22:a:2305:U:H4'	27:f:133:ARG:NH1	2.33	0.42
53:V:24:G:H2'	53:V:25:A:H8	1.84	0.42
1:A:71:A:C4	1:A:72:A:C8	3.07	0.42
1:A:834:U:H2'	1:A:835:U:C6	2.54	0.42
1:A:1268:G:H2'	1:A:1269:A:C8	2.54	0.42
22:a:265:A:N1	22:a:427:U:O2'	2.47	0.42
22:a:1438:U:H2'	22:a:1439:A:H8	1.84	0.42
22:a:1473:G:H2'	22:a:1474:U:H6	1.84	0.42
22:a:1509:A:O2'	22:a:1510:G:H8	2.03	0.42
22:a:1581:G:H2'	22:a:1582:C:C6	2.54	0.42
22:a:2747:G:O6	22:a:2755:C:H5''	2.19	0.42
22:a:2848:G:HO2'	22:a:2867:G:H22	1.65	0.42
22:a:2885:G:O6	47:z:40:ARG:NH1	2.52	0.42
1:A:62:U:OP1	1:A:385:C:O2'	2.35	0.42
1:A:470:C:H2'	1:A:471:U:H6	1.82	0.42
9:I:41:ARG:HD2	9:I:41:ARG:HA	1.70	0.42
13:M:88:GLY:O	13:M:92:ARG:HG3	2.19	0.42
21:U:7:ARG:NE	21:U:10:GLU:OE2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:a:284:U:H2'	22:a:285:G:H5''	2.02	0.42
22:a:641:U:O2'	22:a:2350:C:OP1	2.35	0.42
22:a:852:U:H2'	22:a:853:C:H6	1.84	0.42
53:V:71:A:H2'	53:V:72:C:C6	2.53	0.42
1:A:1071:C:H2'	1:A:1072:G:C8	2.53	0.42
1:A:1225:A:H2'	1:A:1226:C:C6	2.55	0.42
3:C:45:LYS:HE3	3:C:45:LYS:HB3	1.83	0.42
3:C:180:ALA:HA	3:C:206:GLU:HA	2.00	0.42
21:U:11:PRO:HG2	21:U:14:VAL:HG13	2.02	0.42
22:a:38:A:H2'	22:a:39:G:O4'	2.20	0.42
22:a:221:A:N1	22:a:265:A:O2'	2.49	0.42
22:a:441:U:H2'	22:a:442:G:C8	2.54	0.42
22:a:468:G:N7	49:1:39:ARG:NH2	2.51	0.42
22:a:639:U:C2	22:a:640:C:C5	3.07	0.42
22:a:857:G:H2'	22:a:858:G:O4'	2.19	0.42
22:a:1115:G:O2'	22:a:1116:G:H8	2.02	0.42
22:a:1183:U:H2'	22:a:1184:U:C6	2.54	0.42
22:a:1869:G:N2	22:a:1871:A:H3'	2.34	0.42
22:a:2812:G:H2'	22:a:2813:A:C8	2.55	0.42
1:A:8:A:C6	4:D:206:LYS:HB3	2.54	0.42
1:A:166:U:H2'	1:A:167:A:C8	2.55	0.42
1:A:270:A:H2'	1:A:271:C:C6	2.55	0.42
10:J:15:HIS:O	10:J:18:ILE:HG22	2.18	0.42
13:M:101:ARG:NE	13:M:104:THR:HG1	2.14	0.42
22:a:2072:C:H2'	22:a:2073:C:H6	1.84	0.42
22:a:2795:C:H2'	22:a:2796:U:H6	1.85	0.42
26:e:91:ASP:OD1	26:e:91:ASP:N	2.53	0.42
27:f:46:ASP:OD1	27:f:48:LYS:NZ	2.42	0.42
28:g:38:ASN:OD1	28:g:39:ASP:N	2.52	0.42
40:s:19:LYS:HA	40:s:22:THR:HG22	2.00	0.42
53:V:44:G:H2'	53:V:45:U:C6	2.54	0.42
1:A:8:A:N6	4:D:202:GLU:O	2.52	0.42
1:A:235:C:H2'	1:A:236:A:H8	1.83	0.42
1:A:335:C:H2'	1:A:336:A:C8	2.53	0.42
1:A:1378:C:OP1	7:G:5:ARG:NH1	2.53	0.42
12:L:110:ARG:HB3	12:L:119:VAL:HG11	2.01	0.42
22:a:1494:A:H2'	22:a:1495:A:C8	2.55	0.42
22:a:2650:U:H2'	22:a:2651:C:C6	2.54	0.42
28:g:29:LYS:HE2	28:g:80:THR:O	2.20	0.42
37:p:97:ASP:OD2	38:q:13:ARG:NE	2.51	0.42
42:u:64:VAL:HG12	42:u:69:GLU:HG2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:A:H1'	1:A:613:C:O2'	2.20	0.42
1:A:920:U:H2'	1:A:921:U:C6	2.53	0.42
1:A:1001:C:H2'	1:A:1002:G:C8	2.55	0.42
7:G:54:SER:HB3	7:G:56:LYS:HG2	2.01	0.42
15:O:10:LYS:O	15:O:14:GLU:HG3	2.19	0.42
20:T:43:ASP:OD2	20:T:46:ALA:N	2.41	0.42
22:a:1269:A:H2'	22:a:1270:C:C6	2.55	0.42
22:a:1975:G:H2'	22:a:1976:U:O4'	2.20	0.42
22:a:2219:U:H2'	22:a:2220:U:O4'	2.20	0.42
26:e:120:VAL:HA	26:e:188:MET:O	2.19	0.42
30:i:69:ARG:O	30:i:89:PHE:HB3	2.19	0.42
1:A:67:C:H2'	1:A:68:G:C8	2.55	0.42
1:A:1120:C:H2'	1:A:1121:U:C6	2.54	0.42
1:A:1250:A:H2'	1:A:1251:A:C8	2.55	0.42
1:A:1394:A:N1	1:A:1500:A:O2'	2.50	0.42
3:C:46:GLU:HG2	3:C:87:LEU:HD21	2.02	0.42
10:J:5:ARG:HH12	10:J:7:ARG:HE	1.68	0.42
18:R:10:PHE:O	18:R:12:ARG:HD2	2.20	0.42
22:a:475:C:H2'	22:a:476:G:O4'	2.19	0.42
22:a:2405:G:H1'	22:a:2412:A:N6	2.34	0.42
22:a:2502:G:H5''	22:a:2503:2MA:H5''	2.01	0.42
27:f:14:LYS:O	27:f:17:MET:HB2	2.20	0.42
39:r:37:THR:HG23	39:r:38:TYR:CD2	2.55	0.42
13:M:89:LEU:O	13:M:93:ARG:HG3	2.18	0.42
18:R:49:ALA:O	18:R:53:ARG:HG3	2.20	0.42
22:a:566:U:H5''	32:k:29:LYS:HE2	2.01	0.42
22:a:753:A:H2'	22:a:754:U:C6	2.55	0.42
22:a:1187:G:OP1	38:q:85:LYS:NZ	2.41	0.42
22:a:2547:A:H2'	22:a:2548:U:H6	1.83	0.42
22:a:2740:A:H2'	22:a:2741:A:C8	2.54	0.42
22:a:2895:G:H2'	22:a:2896:C:C6	2.55	0.42
26:e:106:LYS:HE2	26:e:200:LEU:HB3	2.00	0.42
26:e:112:LEU:HD13	26:e:186:VAL:HG11	2.02	0.42
27:f:121:SER:HB2	27:f:128:TYR:CE1	2.55	0.42
41:t:81:ASP:OD2	41:t:96:PHE:HB3	2.20	0.42
53:V:20:C:O4'	53:V:22:A:H5'	2.19	0.42
1:A:74:A:H2'	1:A:75:G:C8	2.55	0.42
2:B:7:ARG:O	2:B:11:LYS:HG3	2.20	0.42
18:R:23:TYR:HE2	18:R:65:LEU:HD12	1.85	0.42
22:a:586:A:N1	22:a:809:G:O2'	2.50	0.42
22:a:1447:C:H2'	22:a:1448:G:H8	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:c:227:PRO:HD3	24:c:234:GLY:CA	2.50	0.42
28:g:131:ILE:HG22	28:g:133:LEU:HD22	2.02	0.42
33:l:41:LEU:HG	33:l:96:ILE:HG13	2.00	0.42
44:w:64:ILE:O	44:w:68:LEU:HG	2.20	0.42
52:4:2:LYS:HB2	52:4:5:ILE:HD11	2.02	0.42
1:A:715:A:H2'	1:A:716:A:C8	2.55	0.41
1:A:745:G:H2'	1:A:746:A:C8	2.55	0.41
1:A:904:U:H2'	1:A:905:U:C6	2.55	0.41
1:A:1516:2MG:N1	1:A:1519:MA6:OP2	2.52	0.41
22:a:101:A:N3	22:a:101:A:H2'	2.35	0.41
22:a:1747:U:H2'	22:a:1748:C:C6	2.55	0.41
22:a:2037:A:H2'	22:a:2038:G:C8	2.55	0.41
27:f:38:MET:SD	27:f:53:ALA:HB1	2.60	0.41
28:g:22:GLN:OE1	28:g:55:ARG:NH1	2.52	0.41
34:m:29:VAL:HG11	34:m:75:ILE:HG23	2.02	0.41
35:n:4:LYS:O	35:n:8:ILE:HG12	2.20	0.41
35:n:53:THR:O	35:n:59:ALA:HB2	2.20	0.41
1:A:864:A:H2'	1:A:865:A:C8	2.55	0.41
5:E:83:HIS:CG	8:H:96:MET:HB2	2.55	0.41
7:G:57:SER:OG	7:G:60:GLU:HG3	2.19	0.41
21:U:39:GLU:OE2	21:U:44:GLU:HB3	2.19	0.41
22:a:30:G:H2'	22:a:31:C:C6	2.55	0.41
22:a:635:C:O2'	22:a:639:U:OP1	2.35	0.41
22:a:1410:G:H2'	22:a:1411:U:H6	1.83	0.41
22:a:2097:A:H2'	22:a:2098:U:H6	1.85	0.41
22:a:2501:C:H4'	22:a:2502:G:OP2	2.20	0.41
22:a:2751:G:C4	28:g:3:ARG:HD2	2.55	0.41
25:d:160:LYS:O	25:d:161:MET:HB3	2.20	0.41
27:f:64:LYS:H	52:4:6:HIS:CE1	2.38	0.41
53:V:67:G:H2'	53:V:68:G:C8	2.54	0.41
55:X:5:C:O2'	55:X:6:C:O4'	2.37	0.41
55:X:49:G:H2'	55:X:50:C:C6	2.55	0.41
1:A:499:A:O4'	1:A:547:A:N6	2.53	0.41
7:G:149:LYS:HD2	11:K:61:PHE:CZ	2.56	0.41
22:a:18:U:H2'	22:a:19:A:C8	2.56	0.41
22:a:111:A:H4'	45:x:63:ALA:HB1	2.01	0.41
22:a:1808:A:H3'	22:a:1809:A:C8	2.55	0.41
22:a:2783:U:H2'	22:a:2784:U:C6	2.55	0.41
31:j:120:PRO:HB2	36:o:66:ASN:OD1	2.20	0.41
34:m:69:ARG:C	34:m:71:ARG:H	2.28	0.41
1:A:158:G:H2'	1:A:159:G:C8	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:A:H2'	1:A:264:C:C6	2.56	0.41
22:a:55:G:O2'	22:a:127:A:N1	2.49	0.41
22:a:127:A:N6	49:1:46:LYS:HD2	2.35	0.41
22:a:354:A:H2'	22:a:355:U:O4'	2.21	0.41
22:a:480:A:O2'	41:t:43:LYS:O	2.39	0.41
22:a:1829:A:C8	22:a:1830:C:C5	3.09	0.41
22:a:2300:C:H2'	22:a:2301:C:C6	2.54	0.41
22:a:2680:U:O2'	22:a:2681:C:H5'	2.20	0.41
27:f:140:GLU:OE2	52:4:26:SER:OG	2.25	0.41
40:s:11:LEU:HB2	45:x:26:PHE:HE1	1.84	0.41
51:3:1:MET:HG3	51:3:34:LYS:HB3	2.02	0.41
1:A:456:A:H3'	1:A:457:G:H8	1.85	0.41
1:A:493:A:H2'	1:A:494:G:C8	2.55	0.41
1:A:502:A:H2'	1:A:503:C:O4'	2.20	0.41
1:A:678:U:H2'	1:A:679:C:H6	1.86	0.41
1:A:1119:C:H2'	1:A:1120:C:C6	2.56	0.41
4:D:28:ILE:O	4:D:34:ILE:HD11	2.21	0.41
18:R:26:ILE:CG2	18:R:30:LYS:HE3	2.49	0.41
22:a:58:G:O2'	22:a:73:A:N1	2.51	0.41
22:a:155:A:H2'	22:a:156:A:H8	1.83	0.41
22:a:282:A:N6	22:a:359:G:O6	2.53	0.41
22:a:810:U:C4	32:k:29:LYS:O	2.73	0.41
22:a:1572:A:H2'	22:a:1573:G:C8	2.56	0.41
23:b:57:A:H2'	23:b:58:A:H8	1.86	0.41
27:f:176:PRO:HG3	52:4:44:PHE:CZ	2.56	0.41
34:m:8:ARG:HB2	34:m:43:GLU:HG3	2.01	0.41
43:v:37:ILE:HG21	43:v:80:ILE:HG21	2.03	0.41
1:A:217:C:H2'	1:A:218:U:H6	1.85	0.41
1:A:977:A:O2'	1:A:979:C:OP2	2.35	0.41
1:A:1155:A:H2'	1:A:1156:G:O4'	2.20	0.41
1:A:1163:A:H2'	1:A:1164:G:H8	1.85	0.41
1:A:1338:G:H2'	1:A:1339:A:C8	2.55	0.41
2:B:94:HIS:ND1	2:B:146:ASN:O	2.54	0.41
22:a:1152:C:H4'	37:p:77:SER:HA	2.01	0.41
22:a:1231:U:H2'	22:a:1232:G:H8	1.85	0.41
22:a:1292:G:H2'	22:a:1293:C:H6	1.85	0.41
22:a:1523:U:O2'	22:a:1524:G:OP1	2.36	0.41
22:a:2308:G:H2'	22:a:2308:G:N3	2.36	0.41
22:a:2677:G:H2'	22:a:2678:C:C6	2.56	0.41
31:j:58:LEU:HD11	31:j:86:LEU:HG	2.02	0.41
39:r:25:ARG:NH2	39:r:74:ILE:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:r:37:THR:HB	39:r:48:LYS:NZ	2.35	0.41
1:A:382:A:H2'	1:A:383:A:C8	2.56	0.41
13:M:86:TYR:O	13:M:90:ARG:HG2	2.21	0.41
16:P:18:GLN:HA	16:P:38:PHE:HA	2.03	0.41
21:U:51:SER:O	21:U:54:LYS:HG2	2.20	0.41
22:a:191:A:H2'	22:a:192:C:H6	1.81	0.41
22:a:2552:OMU:HM23	22:a:2554:U:C6	2.55	0.41
27:f:58:ALA:O	27:f:62:GLY:N	2.53	0.41
30:i:13:ARG:HH22	30:i:49:ASP:HB3	1.86	0.41
31:j:1:MET:HE3	31:j:32:TYR:CZ	2.56	0.41
31:j:42:THR:HA	31:j:56:ASP:O	2.20	0.41
1:A:616:G:O2'	16:P:47:GLU:HG3	2.21	0.41
1:A:1384:C:H2'	1:A:1385:G:C8	2.56	0.41
11:K:28:ASN:O	11:K:57:LYS:HE3	2.20	0.41
22:a:2:G:H2'	22:a:3:U:H6	1.86	0.41
22:a:288:U:H2'	22:a:289:G:H8	1.86	0.41
22:a:565:C:H2'	22:a:566:U:O4'	2.21	0.41
22:a:846:U:H4'	22:a:847:U:C5	2.49	0.41
22:a:1842:G:H2'	22:a:1843:C:C6	2.55	0.41
40:s:28:ASN:HD22	40:s:89:GLU:HA	1.86	0.41
1:A:34:C:H2'	1:A:35:G:C8	2.56	0.41
1:A:185:U:H2'	1:A:186:C:H6	1.86	0.41
1:A:1130:A:H2'	1:A:1131:G:H8	1.86	0.41
2:B:43:LEU:HD12	2:B:46:THR:OG1	2.21	0.41
5:E:82:GLN:HG2	5:E:147:MET:HE3	2.02	0.41
8:H:11:LEU:HD22	8:H:75:ILE:HD11	2.03	0.41
8:H:22:LYS:O	8:H:65:TYR:OH	2.32	0.41
18:R:13:PHE:HD1	18:R:18:VAL:HG21	1.86	0.41
22:a:151:C:H2'	22:a:152:A:C8	2.56	0.41
22:a:417:C:H2'	22:a:418:C:C6	2.56	0.41
22:a:1357:C:H2'	22:a:1358:G:O4'	2.21	0.41
22:a:1394:U:H2'	22:a:1395:A:O4'	2.21	0.41
22:a:1437:C:H2'	22:a:1438:U:C6	2.56	0.41
22:a:1587:G:H2'	22:a:1588:G:H8	1.86	0.41
22:a:1842:G:H2'	22:a:1843:C:H6	1.85	0.41
22:a:1874:C:H2'	22:a:1875:G:O4'	2.20	0.41
22:a:2241:A:H2'	22:a:2242:G:C8	2.56	0.41
22:a:2469:A:H2'	22:a:2470:G:O4'	2.21	0.41
28:g:155:GLU:OE2	28:g:158:LYS:N	2.51	0.41
37:p:109:LEU:HD23	38:q:49:ILE:HD13	2.03	0.41
41:t:86:ARG:NE	41:t:88:GLU:OE2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:V:25:A:H2'	53:V:26:C:O4'	2.20	0.41
55:X:49:G:H2'	55:X:50:C:H6	1.86	0.41
55:X:51:U:H2'	55:X:52:G:C8	2.55	0.41
1:A:1118:U:H2'	1:A:1119:C:H6	1.86	0.41
1:A:1228:C:OP2	13:M:110:LYS:NZ	2.48	0.41
1:A:1414:U:H2'	1:A:1415:G:H8	1.86	0.41
1:A:1494:G:H4'	22:a:1913:A:C8	2.56	0.41
2:B:70:VAL:HB	2:B:163:VAL:HG22	2.03	0.41
2:B:97:LEU:HB2	2:B:100:MET:HG3	2.02	0.41
3:C:77:ILE:O	3:C:81:GLY:HA2	2.21	0.41
5:E:110:ALA:O	5:E:114:VAL:HG23	2.21	0.41
6:F:10:VAL:HG21	6:F:18:VAL:HG22	2.02	0.41
13:M:74:SER:OG	13:M:78:LYS:NZ	2.54	0.41
19:S:50:ALA:HA	19:S:58:VAL:O	2.21	0.41
22:a:1957:C:H2'	22:a:1958:C:C6	2.56	0.41
22:a:2590:A:H2'	22:a:2591:C:H6	1.86	0.41
25:d:172:VAL:HG23	25:d:194:PRO:HD3	2.03	0.41
30:i:18:VAL:HG22	30:i:140:LEU:HB3	2.02	0.41
31:j:18:ARG:HB3	31:j:45:GLU:HB2	2.02	0.41
1:A:32:A:H2'	1:A:33:A:C8	2.56	0.40
1:A:477:C:H2'	1:A:478:A:H8	1.83	0.40
1:A:579:A:O2'	15:O:54:ARG:NH1	2.39	0.40
1:A:1477:U:H2'	1:A:1478:U:C6	2.56	0.40
4:D:33:LYS:HA	4:D:33:LYS:HD2	1.80	0.40
4:D:85:ASN:HB3	4:D:88:GLU:CD	2.46	0.40
22:a:184:C:H2'	22:a:185:G:C8	2.56	0.40
22:a:215:G:H4'	22:a:216:A:H4'	2.02	0.40
44:w:37:ARG:HG2	44:w:48:THR:OG1	2.20	0.40
52:4:37:CYS:O	52:4:41:HIS:HB2	2.22	0.40
1:A:780:A:H5''	11:K:125:LYS:HD3	2.02	0.40
22:a:1001:A:H2'	22:a:1002:G:O4'	2.21	0.40
22:a:1223:G:N7	38:q:71:LYS:NZ	2.69	0.40
22:a:1239:G:H2'	22:a:1240:U:O4'	2.20	0.40
22:a:1636:U:H2'	22:a:1637:A:C8	2.56	0.40
22:a:1769:U:O2'	22:a:1958:C:OP1	2.37	0.40
22:a:2030:6MZ:C2	22:a:2499:C:H5''	2.51	0.40
27:f:8:TYR:HB2	27:f:173:PHE:HZ	1.85	0.40
27:f:13:VAL:O	27:f:17:MET:HG2	2.21	0.40
35:n:33:ARG:HG2	35:n:34:HIS:CD2	2.56	0.40
39:r:17:VAL:HG11	39:r:103:ILE:HG12	2.03	0.40
40:s:22:THR:O	40:s:26:LYS:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:29:ILE:HD13	6:F:64:VAL:HG11	2.04	0.40
20:T:51:PHE:C	20:T:51:PHE:CD1	3.00	0.40
22:a:285:G:C5	22:a:356:G:C6	3.10	0.40
22:a:454:A:H4'	22:a:455:C:OP2	2.21	0.40
22:a:1153:C:H2'	22:a:1154:G:O4'	2.22	0.40
22:a:1599:U:H2'	22:a:1600:C:C6	2.56	0.40
23:b:91:C:H2'	23:b:92:C:H6	1.86	0.40
27:f:40:VAL:O	27:f:40:VAL:HG12	2.22	0.40
42:u:26:PHE:CZ	42:u:42:LEU:HD12	2.56	0.40
1:A:196:A:N3	1:A:222:C:H1'	2.37	0.40
1:A:865:A:H5'	1:A:1078:U:O4	2.22	0.40
6:F:18:VAL:HG11	6:F:58:HIS:CE1	2.56	0.40
6:F:25:TYR:HD2	6:F:62:MET:HE1	1.85	0.40
11:K:114:THR:HG21	21:U:32:VAL:HG11	2.04	0.40
22:a:286:U:H2'	22:a:287:G:C8	2.56	0.40
22:a:957:C:H5'	33:l:75:GLU:OE1	2.21	0.40
22:a:1169:A:H2'	22:a:1170:C:C6	2.56	0.40
22:a:1721:G:O2'	22:a:1739:A:N6	2.52	0.40
22:a:1807:G:H5'	22:a:1808:A:OP2	2.21	0.40
22:a:2455:G:H2'	22:a:2456:C:H6	1.86	0.40
22:a:2514:U:H2'	22:a:2515:C:H6	1.84	0.40
22:a:2661:G:H8	22:a:2661:G:OP2	2.03	0.40
39:r:29:VAL:HB	39:r:55:ILE:HD11	2.04	0.40
1:A:321:A:H2'	1:A:322:C:C6	2.57	0.40
1:A:678:U:H2'	1:A:679:C:C6	2.57	0.40
1:A:1003:G:N2	1:A:1005:A:H5'	2.37	0.40
1:A:1007:U:H2'	1:A:1008:U:H6	1.85	0.40
2:B:20:THR:HG22	2:B:39:HIS:CE1	2.57	0.40
4:D:107:PHE:CG	4:D:145:ILE:HD11	2.56	0.40
22:a:208:C:H2'	22:a:209:C:H6	1.85	0.40
22:a:747:5MU:O2	22:a:2014:A:H1'	2.21	0.40
22:a:993:G:OP2	37:p:51:ARG:NH2	2.54	0.40
22:a:1447:C:H2'	22:a:1448:G:C8	2.57	0.40
22:a:1496:A:N3	22:a:1577:C:O2'	2.51	0.40
22:a:2373:G:H2'	22:a:2374:C:C6	2.57	0.40
22:a:2749:A:H5''	28:g:2:SER:HB3	2.04	0.40
26:e:189:THR:O	26:e:193:VAL:HG23	2.22	0.40
30:i:35:ARG:HB3	30:i:54:ILE:HD11	2.04	0.40
37:p:78:LYS:O	37:p:117:LEU:HD11	2.21	0.40
53:V:49:A:H2'	53:V:50:G:C8	2.56	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	
2	B	222/241 (92%)	202 (91%)	20 (9%)	0	100	100
3	C	204/233 (88%)	193 (95%)	11 (5%)	0	100	100
4	D	203/206 (98%)	199 (98%)	4 (2%)	0	100	100
5	E	154/167 (92%)	150 (97%)	4 (3%)	0	100	100
6	F	101/135 (75%)	95 (94%)	6 (6%)	0	100	100
7	G	151/179 (84%)	144 (95%)	7 (5%)	0	100	100
8	H	127/130 (98%)	121 (95%)	6 (5%)	0	100	100
9	I	125/130 (96%)	120 (96%)	5 (4%)	0	100	100
10	J	96/103 (93%)	89 (93%)	6 (6%)	1 (1%)	12	20
11	K	113/129 (88%)	107 (95%)	6 (5%)	0	100	100
12	L	120/124 (97%)	114 (95%)	6 (5%)	0	100	100
13	M	113/118 (96%)	102 (90%)	11 (10%)	0	100	100
14	N	98/101 (97%)	93 (95%)	5 (5%)	0	100	100
15	O	86/89 (97%)	81 (94%)	5 (6%)	0	100	100
16	P	79/82 (96%)	71 (90%)	8 (10%)	0	100	100
17	Q	77/84 (92%)	76 (99%)	1 (1%)	0	100	100
18	R	64/75 (85%)	61 (95%)	3 (5%)	0	100	100
19	S	82/92 (89%)	79 (96%)	3 (4%)	0	100	100
20	T	84/87 (97%)	83 (99%)	1 (1%)	0	100	100
21	U	68/71 (96%)	68 (100%)	0	0	100	100
24	c	269/273 (98%)	263 (98%)	6 (2%)	0	100	100
25	d	206/209 (99%)	198 (96%)	7 (3%)	1 (0%)	24	37
26	e	199/201 (99%)	193 (97%)	6 (3%)	0	100	100
27	f	175/179 (98%)	163 (93%)	12 (7%)	0	100	100
28	g	174/177 (98%)	156 (90%)	17 (10%)	1 (1%)	21	32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
29	h	39/149 (26%)	35 (90%)	4 (10%)	0	100	100
30	i	140/142 (99%)	139 (99%)	1 (1%)	0	100	100
31	j	121/123 (98%)	118 (98%)	3 (2%)	0	100	100
32	k	142/144 (99%)	134 (94%)	8 (6%)	0	100	100
33	l	132/136 (97%)	129 (98%)	3 (2%)	0	100	100
34	m	116/127 (91%)	112 (97%)	4 (3%)	0	100	100
35	n	114/117 (97%)	111 (97%)	3 (3%)	0	100	100
36	o	112/115 (97%)	110 (98%)	2 (2%)	0	100	100
37	p	115/118 (98%)	115 (100%)	0	0	100	100
38	q	101/103 (98%)	98 (97%)	3 (3%)	0	100	100
39	r	108/110 (98%)	106 (98%)	2 (2%)	0	100	100
40	s	91/100 (91%)	82 (90%)	9 (10%)	0	100	100
41	t	100/104 (96%)	90 (90%)	10 (10%)	0	100	100
42	u	92/94 (98%)	90 (98%)	2 (2%)	0	100	100
43	v	76/85 (89%)	72 (95%)	4 (5%)	0	100	100
44	w	75/78 (96%)	75 (100%)	0	0	100	100
45	x	60/63 (95%)	56 (93%)	4 (7%)	0	100	100
46	y	56/59 (95%)	53 (95%)	3 (5%)	0	100	100
47	z	54/57 (95%)	52 (96%)	2 (4%)	0	100	100
48	0	49/55 (89%)	49 (100%)	0	0	100	100
49	1	44/46 (96%)	44 (100%)	0	0	100	100
50	2	62/65 (95%)	58 (94%)	3 (5%)	1 (2%)	7	11
51	3	36/38 (95%)	36 (100%)	0	0	100	100
52	4	56/70 (80%)	52 (93%)	4 (7%)	0	100	100
All	All	5481/5913 (93%)	5237 (96%)	240 (4%)	4 (0%)	49	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
25	d	149	ASN
10	J	57	VAL
50	2	32	ILE
28	g	47	ASP

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	
2	B	186/199 (94%)	186 (100%)	0	100	100
3	C	170/190 (90%)	170 (100%)	0	100	100
4	D	172/173 (99%)	172 (100%)	0	100	100
5	E	119/126 (94%)	119 (100%)	0	100	100
6	F	90/116 (78%)	90 (100%)	0	100	100
7	G	126/147 (86%)	126 (100%)	0	100	100
8	H	104/105 (99%)	104 (100%)	0	100	100
9	I	105/107 (98%)	105 (100%)	0	100	100
10	J	86/90 (96%)	86 (100%)	0	100	100
11	K	89/98 (91%)	89 (100%)	0	100	100
12	L	102/103 (99%)	102 (100%)	0	100	100
13	M	93/96 (97%)	93 (100%)	0	100	100
14	N	83/84 (99%)	83 (100%)	0	100	100
15	O	76/77 (99%)	76 (100%)	0	100	100
16	P	65/65 (100%)	65 (100%)	0	100	100
17	Q	73/78 (94%)	73 (100%)	0	100	100
18	R	57/65 (88%)	57 (100%)	0	100	100
19	S	72/79 (91%)	72 (100%)	0	100	100
20	T	65/66 (98%)	65 (100%)	0	100	100
21	U	60/61 (98%)	60 (100%)	0	100	100
24	c	216/218 (99%)	216 (100%)	0	100	100
25	d	163/163 (100%)	163 (100%)	0	100	100
26	e	165/165 (100%)	165 (100%)	0	100	100
27	f	148/150 (99%)	148 (100%)	0	100	100
28	g	137/138 (99%)	137 (100%)	0	100	100
29	h	32/114 (28%)	32 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
30	i	116/116 (100%)	116 (100%)	0	100	100
31	j	104/104 (100%)	104 (100%)	0	100	100
32	k	103/103 (100%)	103 (100%)	0	100	100
33	l	107/107 (100%)	107 (100%)	0	100	100
34	m	98/103 (95%)	98 (100%)	0	100	100
35	n	86/87 (99%)	86 (100%)	0	100	100
36	o	99/100 (99%)	99 (100%)	0	100	100
37	p	89/90 (99%)	89 (100%)	0	100	100
38	q	84/84 (100%)	84 (100%)	0	100	100
39	r	93/93 (100%)	93 (100%)	0	100	100
40	s	80/84 (95%)	80 (100%)	0	100	100
41	t	83/85 (98%)	83 (100%)	0	100	100
42	u	78/78 (100%)	78 (100%)	0	100	100
43	v	58/63 (92%)	58 (100%)	0	100	100
44	w	67/68 (98%)	67 (100%)	0	100	100
45	x	54/55 (98%)	54 (100%)	0	100	100
46	y	48/49 (98%)	48 (100%)	0	100	100
47	z	47/48 (98%)	47 (100%)	0	100	100
48	0	46/49 (94%)	46 (100%)	0	100	100
49	1	38/38 (100%)	38 (100%)	0	100	100
50	2	51/52 (98%)	51 (100%)	0	100	100
51	3	34/34 (100%)	34 (100%)	0	100	100
52	4	55/62 (89%)	55 (100%)	0	100	100
All	All	4572/4825 (95%)	4572 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	51	ASN
2	B	109	GLN
2	B	120	GLN
2	B	190	ASN

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Mol	Chain	Res	Type
3	C	190	HIS
4	D	196	ASN
5	E	77	ASN
5	E	148	ASN
6	F	68	GLN
8	H	4	GLN
8	H	38	ASN
9	I	75	GLN
10	J	58	ASN
11	K	109	ASN
13	M	14	HIS
13	M	105	ASN
20	T	20	HIS
20	T	78	ASN
24	c	37	ASN
24	c	117	GLN
24	c	226	ASN
25	d	126	ASN
26	e	115	GLN
27	f	5	HIS
27	f	37	ASN
28	g	45	HIS
29	h	18	GLN
34	m	9	GLN
35	n	98	GLN
35	n	104	GLN
37	p	52	GLN
37	p	71	GLN
39	r	9	HIS
46	y	9	GLN
52	4	61	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1516/1542 (98%)	212 (13%)	3 (0%)
22	a	2749/2904 (94%)	298 (10%)	0
23	b	118/120 (98%)	13 (11%)	0
53	V	75/76 (98%)	14 (18%)	0
54	W	10/48 (20%)	1 (10%)	0
55	X	75/78 (96%)	29 (38%)	3 (4%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	4543/4768 (95%)	567 (12%)	6 (0%)

All (567) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	5	U
1	A	6	G
1	A	7	A
1	A	9	G
1	A	22	G
1	A	32	A
1	A	39	G
1	A	47	C
1	A	48	C
1	A	50	A
1	A	51	A
1	A	60	A
1	A	64	G
1	A	69	G
1	A	71	A
1	A	72	A
1	A	74	A
1	A	80	A
1	A	81	A
1	A	83	C
1	A	84	U
1	A	86	G
1	A	87	C
1	A	88	U
1	A	90	C
1	A	91	U
1	A	94	G
1	A	95	C
1	A	116	A
1	A	121	U
1	A	122	G
1	A	131	A
1	A	141	G
1	A	144	G
1	A	148	G
1	A	163	C
1	A	175	C

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Mol	Chain	Res	Type
1	A	182	A
1	A	191	G
1	A	195	A
1	A	197	A
1	A	200	G
1	A	204	G
1	A	226	G
1	A	236	A
1	A	245	U
1	A	247	G
1	A	251	G
1	A	262	A
1	A	266	G
1	A	267	C
1	A	280	C
1	A	289	G
1	A	321	A
1	A	328	C
1	A	338	A
1	A	347	G
1	A	351	G
1	A	352	C
1	A	354	G
1	A	367	U
1	A	372	C
1	A	373	A
1	A	380	G
1	A	384	G
1	A	397	A
1	A	406	G
1	A	411	A
1	A	412	A
1	A	413	G
1	A	414	A
1	A	421	U
1	A	422	C
1	A	423	G
1	A	424	G
1	A	429	U
1	A	446	G
1	A	453	G
1	A	457	G

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Mol	Chain	Res	Type
1	A	458	U
1	A	465	A
1	A	467	U
1	A	468	A
1	A	481	G
1	A	484	G
1	A	486	U
1	A	495	A
1	A	496	A
1	A	497	G
1	A	499	A
1	A	509	A
1	A	511	C
1	A	518	C
1	A	520	A
1	A	521	G
1	A	529	G
1	A	531	U
1	A	532	A
1	A	547	A
1	A	559	A
1	A	564	C
1	A	572	A
1	A	573	A
1	A	576	C
1	A	579	A
1	A	633	G
1	A	637	C
1	A	639	G
1	A	650	G
1	A	653	U
1	A	665	A
1	A	721	G
1	A	723	U
1	A	724	G
1	A	747	A
1	A	755	G
1	A	777	A
1	A	793	U
1	A	794	A
1	A	802	A
1	A	815	A

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Mol	Chain	Res	Type
1	A	817	C
1	A	821	G
1	A	847	G
1	A	890	G
1	A	914	A
1	A	926	G
1	A	934	C
1	A	942	G
1	A	960	U
1	A	961	U
1	A	966	2MG
1	A	969	A
1	A	975	A
1	A	976	G
1	A	977	A
1	A	992	U
1	A	993	G
1	A	1001	C
1	A	1002	G
1	A	1003	G
1	A	1004	A
1	A	1009	U
1	A	1020	G
1	A	1022	A
1	A	1027	C
1	A	1030	U
1	A	1031	C
1	A	1036	A
1	A	1037	C
1	A	1040	U
1	A	1053	G
1	A	1065	U
1	A	1070	U
1	A	1086	U
1	A	1094	G
1	A	1095	U
1	A	1101	A
1	A	1130	A
1	A	1137	C
1	A	1139	G
1	A	1159	U
1	A	1160	G

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Mol	Chain	Res	Type
1	A	1168	U
1	A	1184	G
1	A	1196	A
1	A	1197	A
1	A	1213	A
1	A	1214	C
1	A	1227	A
1	A	1236	A
1	A	1238	A
1	A	1257	A
1	A	1258	G
1	A	1260	G
1	A	1280	A
1	A	1285	A
1	A	1287	A
1	A	1298	U
1	A	1300	G
1	A	1301	U
1	A	1302	C
1	A	1305	G
1	A	1317	C
1	A	1320	C
1	A	1338	G
1	A	1346	A
1	A	1353	G
1	A	1363	A
1	A	1364	U
1	A	1370	G
1	A	1378	C
1	A	1379	G
1	A	1381	U
1	A	1397	C
1	A	1419	G
1	A	1422	G
1	A	1441	A
1	A	1446	A
1	A	1450	U
1	A	1451	U
1	A	1452	C
1	A	1453	G
1	A	1487	G
1	A	1492	A

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Mol	Chain	Res	Type
1	A	1497	G
1	A	1499	A
1	A	1503	A
1	A	1506	U
1	A	1517	G
1	A	1529	G
1	A	1530	G
22	a	10	A
22	a	34	U
22	a	45	G
22	a	51	G
22	a	58	G
22	a	71	A
22	a	74	A
22	a	75	G
22	a	84	A
22	a	101	A
22	a	102	U
22	a	103	A
22	a	114	U
22	a	118	A
22	a	119	A
22	a	120	U
22	a	139	U
22	a	141	G
22	a	142	A
22	a	146	A
22	a	163	C
22	a	165	A
22	a	181	A
22	a	196	A
22	a	199	A
22	a	216	A
22	a	221	A
22	a	222	A
22	a	223	A
22	a	233	A
22	a	248	G
22	a	265	A
22	a	272	A
22	a	277	G
22	a	278	A

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Mol	Chain	Res	Type
22	a	279	A
22	a	285	G
22	a	311	A
22	a	329	G
22	a	330	A
22	a	345	A
22	a	356	G
22	a	361	G
22	a	362	A
22	a	363	G
22	a	386	G
22	a	396	G
22	a	405	U
22	a	411	G
22	a	473	G
22	a	479	A
22	a	481	G
22	a	489	G
22	a	491	G
22	a	504	A
22	a	505	A
22	a	508	A
22	a	509	C
22	a	529	A
22	a	530	G
22	a	531	C
22	a	532	A
22	a	533	G
22	a	545	U
22	a	546	U
22	a	548	G
22	a	549	G
22	a	563	A
22	a	573	U
22	a	575	A
22	a	603	A
22	a	615	U
22	a	627	A
22	a	637	A
22	a	645	C
22	a	646	U
22	a	647	G

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Mol	Chain	Res	Type
22	a	654	A
22	a	655	A
22	a	686	U
22	a	730	A
22	a	738	G
22	a	746	PSU
22	a	747	5MU
22	a	765	C
22	a	775	G
22	a	776	G
22	a	782	A
22	a	784	G
22	a	785	G
22	a	805	G
22	a	812	C
22	a	827	U
22	a	828	U
22	a	845	A
22	a	846	U
22	a	847	U
22	a	859	G
22	a	884	U
22	a	888	C
22	a	890	C
22	a	891	G
22	a	893	C
22	a	895	U
22	a	896	A
22	a	899	A
22	a	910	A
22	a	914	G
22	a	931	U
22	a	932	U
22	a	946	C
22	a	958	U
22	a	961	C
22	a	974	G
22	a	983	A
22	a	996	A
22	a	1012	U
22	a	1013	C
22	a	1033	U

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Mol	Chain	Res	Type
22	a	1040	A
22	a	1045	C
22	a	1046	A
22	a	1047	G
22	a	1108	U
22	a	1111	A
22	a	1112	G
22	a	1128	G
22	a	1132	U
22	a	1133	A
22	a	1135	C
22	a	1142	A
22	a	1144	A
22	a	1206	G
22	a	1253	A
22	a	1256	G
22	a	1271	G
22	a	1272	A
22	a	1273	U
22	a	1300	G
22	a	1301	A
22	a	1329	U
22	a	1345	C
22	a	1352	U
22	a	1365	A
22	a	1379	U
22	a	1383	A
22	a	1414	C
22	a	1416	G
22	a	1417	C
22	a	1421	G
22	a	1428	C
22	a	1452	G
22	a	1453	A
22	a	1460	U
22	a	1476	U
22	a	1482	G
22	a	1493	C
22	a	1508	A
22	a	1509	A
22	a	1510	G
22	a	1515	A

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Mol	Chain	Res	Type
22	a	1524	G
22	a	1531	C
22	a	1535	A
22	a	1536	C
22	a	1537	G
22	a	1538	G
22	a	1566	A
22	a	1569	A
22	a	1578	U
22	a	1583	A
22	a	1584	U
22	a	1585	C
22	a	1592	C
22	a	1607	C
22	a	1608	A
22	a	1647	U
22	a	1648	U
22	a	1649	G
22	a	1674	G
22	a	1715	G
22	a	1729	U
22	a	1730	C
22	a	1731	G
22	a	1738	G
22	a	1764	C
22	a	1773	A
22	a	1782	U
22	a	1800	C
22	a	1801	A
22	a	1808	A
22	a	1816	C
22	a	1829	A
22	a	1848	A
22	a	1858	A
22	a	1868	C
22	a	1870	C
22	a	1871	A
22	a	1884	G
22	a	1906	G
22	a	1929	G
22	a	1930	G
22	a	1936	A

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Mol	Chain	Res	Type
22	a	1937	A
22	a	1955	U
22	a	1964	G
22	a	1967	C
22	a	1970	A
22	a	1971	U
22	a	1972	G
22	a	1991	U
22	a	1993	U
22	a	2023	C
22	a	2031	A
22	a	2033	A
22	a	2043	C
22	a	2055	C
22	a	2056	G
22	a	2059	A
22	a	2060	A
22	a	2061	G
22	a	2062	A
22	a	2069	G7M
22	a	2093	G
22	a	2198	A
22	a	2203	U
22	a	2204	G
22	a	2211	A
22	a	2225	A
22	a	2238	G
22	a	2239	G
22	a	2268	A
22	a	2279	G
22	a	2283	C
22	a	2287	A
22	a	2288	A
22	a	2305	U
22	a	2308	G
22	a	2322	A
22	a	2325	G
22	a	2333	A
22	a	2336	A
22	a	2347	C
22	a	2350	C
22	a	2361	G

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Mol	Chain	Res	Type
22	a	2383	G
22	a	2385	C
22	a	2396	G
22	a	2402	U
22	a	2406	A
22	a	2424	C
22	a	2425	A
22	a	2429	G
22	a	2430	A
22	a	2435	A
22	a	2441	U
22	a	2448	A
22	a	2470	G
22	a	2475	C
22	a	2476	A
22	a	2482	A
22	a	2491	U
22	a	2502	G
22	a	2518	A
22	a	2529	G
22	a	2547	A
22	a	2554	U
22	a	2566	A
22	a	2567	G
22	a	2573	C
22	a	2585	U
22	a	2586	U
22	a	2602	A
22	a	2609	U
22	a	2613	U
22	a	2629	U
22	a	2630	G
22	a	2661	G
22	a	2689	U
22	a	2690	U
22	a	2714	G
22	a	2716	C
22	a	2726	A
22	a	2733	A
22	a	2744	G
22	a	2748	A
22	a	2757	A

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Mol	Chain	Res	Type
22	a	2765	A
22	a	2778	A
22	a	2791	G
22	a	2799	A
22	a	2820	A
22	a	2821	A
22	a	2835	A
22	a	2836	U
22	a	2884	U
22	a	2896	C
22	a	2902	C
23	b	13	G
23	b	24	G
23	b	25	U
23	b	33	G
23	b	35	C
23	b	36	C
23	b	44	G
23	b	56	G
23	b	67	G
23	b	89	U
23	b	90	C
23	b	99	A
23	b	109	A
53	V	5	C
53	V	9	C
53	V	16	A
53	V	17	G
53	V	18	G
53	V	19	C
53	V	20	C
53	V	21	C
53	V	22	A
53	V	23	G
53	V	49	A
53	V	54	5MU
53	V	56	C
53	V	76	A
54	W	35	A
55	X	4	C
55	X	5	C
55	X	6	C

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Mol	Chain	Res	Type
55	X	7	U
55	X	8	U
55	X	9	A
55	X	13	U
55	X	14	A
55	X	15	7SN
55	X	17	C
55	X	17(A)	U
55	X	18	G
55	X	19	G
55	X	20	U
55	X	20(A)	U
55	X	21	A
55	X	22	A
55	X	26	M2G
55	X	27	A
55	X	45	G
55	X	48	5MC
55	X	49	G
55	X	57	7S3
55	X	66	A
55	X	69	G
55	X	70	C
55	X	71	C
55	X	74	C
55	X	75	C

All (6) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	519	C
1	A	1026	G
1	A	1035	A
55	X	6	C
55	X	20	U
55	X	65	C

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

53 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$
22	PSU	a	955	22	18,21,22	3.76	7 (38%)	22,30,33	1.92	5 (22%)
1	2MG	A	1207	1	23,26,27	2.60	7 (30%)	32,38,41	2.14	9 (28%)
1	UR3	A	1498	1	19,22,23	2.86	8 (42%)	26,32,35	1.38	3 (11%)
22	G7M	a	2069	22	23,26,27	2.77	8 (34%)	35,39,42	1.71	9 (25%)
55	5MC	X	48	55	18,22,23	3.76	8 (44%)	26,32,35	1.00	1 (3%)
22	OMU	a	2552	22	19,22,23	2.83	7 (36%)	26,31,34	1.68	5 (19%)
22	5MU	a	747	22	19,22,23	4.69	6 (31%)	28,32,35	3.69	10 (35%)
53	PSU	V	55	53	18,21,22	3.99	6 (33%)	22,30,33	1.90	5 (22%)
1	G7M	A	527	1	23,26,27	2.80	9 (39%)	35,39,42	1.69	8 (22%)
22	6MZ	a	1618	22	22,25,26	2.78	4 (18%)	30,36,39	2.31	9 (30%)
22	OMC	a	2498	56,22	19,22,23	2.95	8 (42%)	26,31,34	0.86	0
55	M2G	X	26	55	24,27,28	1.02	1 (4%)	35,40,43	0.96	2 (5%)
53	PSU	V	13	53	18,21,22	3.89	6 (33%)	22,30,33	1.79	5 (22%)
1	5MC	A	967	1	18,22,23	3.74	8 (44%)	26,32,35	1.00	1 (3%)
22	PSU	a	746	56,22	18,21,22	3.87	7 (38%)	22,30,33	1.65	3 (13%)
25	MEQ	d	150	25	8,9,10	1.03	0	5,10,12	1.44	2 (40%)
33	4D4	l	81	33	9,11,12	1.39	2 (22%)	8,13,15	2.14	4 (50%)
22	2MG	a	2445	22	23,26,27	2.60	8 (34%)	32,38,41	2.21	9 (28%)
22	PSU	a	2604	22	18,21,22	3.86	6 (33%)	22,30,33	1.82	5 (22%)
22	1MG	a	745	22	22,26,27	2.59	7 (31%)	33,39,42	1.66	7 (21%)
55	7S3	X	57	55	22,25,26	2.99	7 (31%)	32,37,40	1.98	9 (28%)
53	5MU	V	54	53	19,22,23	4.71	5 (26%)	28,32,35	3.54	8 (28%)
1	MA6	A	1519	1	23,26,27	1.31	4 (17%)	34,38,41	2.28	11 (32%)
22	3TD	a	1915	22	19,22,23	3.98	6 (31%)	21,32,35	1.58	2 (9%)
55	T6A	X	37	55,56	31,34,35	1.98	7 (22%)	44,49,52	2.14	11 (25%)
1	PSU	A	516	1,56	18,21,22	3.86	6 (33%)	22,30,33	1.77	6 (27%)
55	7SN	X	15	55	25,28,29	2.77	11 (44%)	31,42,45	1.94	6 (19%)
55	OMC	X	56	55	19,22,23	3.07	8 (42%)	26,31,34	0.84	0
53	U8U	V	35	53,54	19,24,25	2.56	6 (31%)	23,34,37	1.16	3 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	IAS	K	119	11	6,7,8	1.09	0	6,8,10	1.40	2 (33%)
22	5MU	a	1939	22	19,22,23	4.73	7 (36%)	28,32,35	3.63	9 (32%)
1	MA6	A	1518	1	23,26,27	1.31	4 (17%)	34,38,41	2.31	12 (35%)
53	2MA	V	38	53	22,25,26	3.76	8 (36%)	33,37,40	2.84	10 (30%)
22	PSU	a	1911	22	18,21,22	3.85	6 (33%)	22,30,33	1.77	5 (22%)
22	6MZ	a	2030	22	22,25,26	2.68	4 (18%)	30,36,39	2.44	11 (36%)
55	PSU	X	55	55	18,21,22	3.95	6 (33%)	22,30,33	1.77	5 (22%)
1	2MG	A	966	1	23,26,27	2.67	7 (30%)	32,38,41	2.26	9 (28%)
22	H2U	a	2449	22	18,21,22	1.11	2 (11%)	21,30,33	0.68	0
1	2MG	A	1516	1	23,26,27	2.64	8 (34%)	32,38,41	2.16	10 (31%)
1	5MC	A	1407	1	18,22,23	3.70	8 (44%)	26,32,35	1.12	3 (11%)
22	5MC	a	1962	22	18,22,23	3.68	8 (44%)	26,32,35	0.98	2 (7%)
22	PSU	a	2580	56,22	18,21,22	3.84	7 (38%)	22,30,33	1.93	6 (27%)
1	4OC	A	1402	1,56	20,23,24	3.08	8 (40%)	26,32,35	0.99	2 (7%)
22	PSU	a	2504	22	18,21,22	3.96	6 (33%)	22,30,33	1.83	4 (18%)
22	2MG	a	1835	22	23,26,27	2.63	7 (30%)	32,38,41	2.14	9 (28%)
12	D2T	L	89	12	7,9,10	2.05	2 (28%)	6,11,13	1.93	3 (50%)
22	OMG	a	2251	56,53,22	23,26,27	2.60	8 (34%)	33,38,41	1.96	9 (27%)
22	PSU	a	1917	22	18,21,22	3.84	6 (33%)	22,30,33	1.84	5 (22%)
22	2MA	a	2503	56,22	22,25,26	3.90	8 (36%)	33,37,40	2.79	9 (27%)
22	PSU	a	2605	22	18,21,22	3.82	6 (33%)	22,30,33	1.86	5 (22%)
33	MS6	l	82	33	5,7,8	1.85	1 (20%)	2,7,9	2.14	1 (50%)
22	PSU	a	2457	22	18,21,22	3.87	6 (33%)	22,30,33	1.93	6 (27%)
55	B8H	X	54	55	20,22,23	3.84	8 (40%)	21,32,35	1.35	1 (4%)

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
22	PSU	a	955	22	-	0/7/25/26	0/2/2/2
1	2MG	A	1207	1	-	0/9/27/28	0/3/3/3
1	UR3	A	1498	1	-	0/7/25/26	0/2/2/2
22	G7M	a	2069	22	-	1/7/25/26	0/3/3/3
55	5MC	X	48	55	-	2/7/25/26	0/2/2/2
22	OMU	a	2552	22	-	1/9/27/28	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	5MU	a	747	22	-	0/7/25/26	0/2/2/2
53	PSU	V	55	53	-	2/7/25/26	0/2/2/2
1	G7M	A	527	1	-	2/7/25/26	0/3/3/3
22	6MZ	a	1618	22	-	0/9/27/28	0/3/3/3
22	OMC	a	2498	56,22	-	0/9/27/28	0/2/2/2
55	M2G	X	26	55	-	2/11/29/30	0/3/3/3
53	PSU	V	13	53	-	2/7/25/26	0/2/2/2
1	5MC	A	967	1	-	0/7/25/26	0/2/2/2
22	PSU	a	746	56,22	-	1/7/25/26	0/2/2/2
25	MEQ	d	150	25	-	5/8/9/11	-
33	4D4	l	81	33	-	1/11/12/14	-
22	2MG	a	2445	22	-	2/9/27/28	0/3/3/3
22	PSU	a	2604	22	-	0/7/25/26	0/2/2/2
22	1MG	a	745	22	-	0/7/25/26	0/3/3/3
55	7S3	X	57	55	-	2/7/25/26	0/3/3/3
53	5MU	V	54	53	-	2/7/25/26	0/2/2/2
1	MA6	A	1519	1	-	2/11/29/30	0/3/3/3
22	3TD	a	1915	22	-	2/7/25/26	0/2/2/2
55	T6A	X	37	55,56	-	6/23/41/42	0/3/3/3
1	PSU	A	516	1,56	-	0/7/25/26	0/2/2/2
55	7SN	X	15	55	-	3/7/29/30	0/3/3/3
55	OMC	X	56	55	-	1/9/27/28	0/2/2/2
53	U8U	V	35	53,54	-	1/9/28/29	0/2/2/2
11	IAS	K	119	11	-	0/7/7/8	-
22	5MU	a	1939	22	-	0/7/25/26	0/2/2/2
1	MA6	A	1518	1	-	0/11/29/30	0/3/3/3
53	2MA	V	38	53	-	0/7/25/26	0/3/3/3
22	PSU	a	1911	22	-	0/7/25/26	0/2/2/2
22	6MZ	a	2030	22	-	2/9/27/28	0/3/3/3
55	PSU	X	55	55	-	1/7/25/26	0/2/2/2
1	2MG	A	966	1	-	2/9/27/28	0/3/3/3
22	H2U	a	2449	22	-	0/7/38/39	0/2/2/2
1	2MG	A	1516	1	-	0/9/27/28	0/3/3/3
1	5MC	A	1407	1	-	0/7/25/26	0/2/2/2
22	5MC	a	1962	22	-	0/7/25/26	0/2/2/2
22	PSU	a	2580	56,22	-	0/7/25/26	0/2/2/2
1	4OC	A	1402	1,56	-	2/9/29/30	0/2/2/2
22	PSU	a	2504	22	-	2/7/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	2MG	a	1835	22	-	0/9/27/28	0/3/3/3
12	D2T	L	89	12	-	2/7/12/14	-
22	OMG	a	2251	56,53,22	-	1/9/27/28	0/3/3/3
22	PSU	a	1917	22	-	0/7/25/26	0/2/2/2
22	2MA	a	2503	56,22	-	1/7/25/26	0/3/3/3
22	PSU	a	2605	22	-	0/7/25/26	0/2/2/2
33	MS6	l	82	33	-	1/4/6/8	-
22	PSU	a	2457	22	-	0/7/25/26	0/2/2/2
55	B8H	X	54	55	-	0/7/25/26	0/2/2/2

All (324) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	1915	3TD	C6-C5	12.25	1.49	1.35
22	a	2503	2MA	C4-N3	12.16	1.49	1.34
22	a	1618	6MZ	C6-N6	11.95	1.47	1.34
22	a	2030	6MZ	C6-N6	11.38	1.46	1.34
53	V	38	2MA	C4-N3	11.32	1.48	1.34
53	V	54	5MU	C6-N1	10.82	1.56	1.38
53	V	54	5MU	C2-N1	10.80	1.55	1.38
22	a	1939	5MU	C6-N1	10.76	1.56	1.38
22	a	747	5MU	C2-N1	10.73	1.55	1.38
22	a	747	5MU	C6-N1	10.67	1.56	1.38
22	a	1939	5MU	C2-N1	10.28	1.54	1.38
53	V	55	PSU	C6-C5	10.03	1.47	1.35
22	a	2504	PSU	C6-C5	9.99	1.46	1.35
55	X	55	PSU	C6-C5	9.97	1.46	1.35
55	X	54	B8H	C2-N3	9.88	1.55	1.38
22	a	2457	PSU	C6-C5	9.69	1.46	1.35
55	X	57	7S3	C2-N3	9.61	1.48	1.30
22	a	1911	PSU	C6-C5	9.56	1.46	1.35
22	a	2580	PSU	C6-C5	9.55	1.46	1.35
1	A	516	PSU	C6-C5	9.55	1.46	1.35
22	a	2604	PSU	C6-C5	9.54	1.46	1.35
22	a	746	PSU	C6-C5	9.53	1.46	1.35
53	V	13	PSU	C6-C5	9.51	1.46	1.35
22	a	2605	PSU	C6-C5	9.50	1.46	1.35
22	a	1917	PSU	C6-C5	9.49	1.46	1.35
1	A	967	5MC	C6-C5	9.45	1.50	1.34
1	A	1407	5MC	C6-C5	9.45	1.50	1.34
22	a	2504	PSU	C2-N1	9.45	1.49	1.36
55	X	48	5MC	C6-C5	9.39	1.50	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	1939	5MU	C4-C5	9.39	1.60	1.44
53	V	55	PSU	C2-N1	9.37	1.49	1.36
22	a	955	PSU	C6-C5	9.36	1.46	1.35
53	V	54	5MU	C4-C5	9.35	1.60	1.44
22	a	1962	5MC	C6-C5	9.33	1.49	1.34
53	V	13	PSU	C2-N1	9.30	1.49	1.36
22	a	747	5MU	C4-C5	9.23	1.60	1.44
22	a	1911	PSU	C2-N1	9.21	1.49	1.36
22	a	2604	PSU	C2-N1	9.21	1.49	1.36
22	a	2457	PSU	C2-N1	9.20	1.49	1.36
55	X	55	PSU	C2-N1	9.18	1.49	1.36
22	a	2580	PSU	C2-N1	9.17	1.49	1.36
22	a	746	PSU	C2-N1	9.16	1.49	1.36
22	a	1917	PSU	C2-N1	9.15	1.49	1.36
1	A	516	PSU	C2-N1	9.11	1.49	1.36
22	a	2605	PSU	C2-N1	9.10	1.49	1.36
22	a	955	PSU	C2-N1	8.80	1.48	1.36
55	X	54	B8H	C2-N1	8.63	1.53	1.38
22	a	1915	3TD	C2-N1	8.54	1.48	1.37
1	A	966	2MG	C2-N3	7.72	1.46	1.31
22	a	1939	5MU	C4-N3	-7.63	1.24	1.38
22	a	1835	2MG	C2-N3	7.61	1.46	1.31
1	A	1516	2MG	C2-N3	7.60	1.46	1.31
1	A	1498	UR3	C2-N1	7.55	1.49	1.38
22	a	2445	2MG	C2-N3	7.47	1.46	1.31
1	A	1207	2MG	C2-N3	7.47	1.46	1.31
22	a	2503	2MA	C2-N3	7.35	1.47	1.34
22	a	2251	OMG	C4-N3	7.23	1.51	1.34
53	V	38	2MA	C2-N3	7.19	1.46	1.34
55	X	48	5MC	C4-N3	7.18	1.46	1.34
53	V	54	5MU	C4-N3	-7.17	1.25	1.38
22	a	747	5MU	C4-N3	-7.15	1.25	1.38
1	A	967	5MC	C4-N3	7.07	1.46	1.34
55	X	55	PSU	C2-N3	6.99	1.49	1.37
55	X	54	B8H	C6-C5	6.99	1.44	1.34
22	a	1962	5MC	C4-N3	6.98	1.45	1.34
53	V	55	PSU	C2-N3	6.94	1.49	1.37
1	A	1407	5MC	C4-N3	6.83	1.45	1.34
53	V	13	PSU	C2-N3	6.77	1.49	1.37
22	a	2504	PSU	C2-N3	6.72	1.49	1.37
22	a	1939	5MU	C6-C5	6.70	1.45	1.34
55	X	57	7S3	C4-N3	6.68	1.49	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1402	4OC	C4-N3	6.68	1.44	1.32
1	A	516	PSU	C2-N3	6.66	1.48	1.37
22	a	1917	PSU	C2-N3	6.64	1.48	1.37
22	a	955	PSU	C2-N3	6.64	1.48	1.37
22	a	1911	PSU	C2-N3	6.62	1.48	1.37
1	A	527	G7M	C4-N3	6.60	1.49	1.34
1	A	1516	2MG	C4-N3	6.60	1.49	1.34
22	a	2552	OMU	C2-N1	6.59	1.49	1.38
22	a	2604	PSU	C2-N3	6.57	1.48	1.37
1	A	966	2MG	C4-N3	6.57	1.49	1.34
22	a	2580	PSU	C2-N3	6.53	1.48	1.37
22	a	2605	PSU	C2-N3	6.50	1.48	1.37
22	a	2457	PSU	C2-N3	6.49	1.48	1.37
55	X	48	5MC	C2-N3	6.47	1.49	1.36
22	a	1835	2MG	C4-N3	6.46	1.49	1.34
22	a	746	PSU	C2-N3	6.45	1.48	1.37
55	X	56	OMC	C2-N3	6.44	1.49	1.36
22	a	2503	2MA	C2-N1	6.43	1.45	1.34
22	a	747	5MU	C6-C5	6.38	1.45	1.34
1	A	527	G7M	C2-N2	6.35	1.49	1.34
22	a	2069	G7M	C4-N3	6.34	1.49	1.34
22	a	2069	G7M	C2-N2	6.32	1.49	1.34
53	V	38	2MA	C2-N1	6.31	1.45	1.34
22	a	1962	5MC	C2-N3	6.29	1.49	1.36
55	X	15	7SN	C4-N3	6.28	1.49	1.34
22	a	2445	2MG	C4-N3	6.27	1.49	1.34
1	A	967	5MC	C2-N3	6.27	1.49	1.36
53	V	54	5MU	C6-C5	6.26	1.44	1.34
22	a	745	1MG	C2-N3	6.23	1.45	1.34
1	A	1207	2MG	C4-N3	6.23	1.49	1.34
22	a	2552	OMU	C2-N3	6.23	1.49	1.38
1	A	1402	4OC	C6-C5	6.20	1.49	1.35
1	A	1498	UR3	C6-C5	6.18	1.49	1.35
22	a	2498	OMC	C2-N3	6.09	1.48	1.36
1	A	1407	5MC	C2-N3	6.04	1.48	1.36
22	a	745	1MG	C4-N3	6.03	1.48	1.34
55	X	56	OMC	C6-C5	6.00	1.49	1.35
22	a	2498	OMC	C6-C5	5.93	1.48	1.35
22	a	745	1MG	C2-N2	5.89	1.44	1.34
22	a	1915	3TD	C6-N1	5.83	1.45	1.36
22	a	2552	OMU	C6-C5	5.82	1.48	1.35
1	A	527	G7M	C2-N3	5.70	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1402	4OC	C2-N3	5.69	1.47	1.36
53	V	35	U8U	C2-N3	5.60	1.50	1.37
22	a	2069	G7M	C2-N3	5.54	1.46	1.33
55	X	37	T6A	C10-N6	5.52	1.49	1.37
53	V	35	U8U	C6-C5	5.50	1.50	1.34
55	X	15	7SN	C2-N3	5.48	1.46	1.33
55	X	54	B8H	C4-N3	5.44	1.48	1.38
55	X	37	T6A	C6-N6	5.43	1.50	1.39
22	a	2251	OMG	C2-N3	5.42	1.46	1.33
53	V	38	2MA	C5-C6	5.33	1.55	1.41
1	A	1498	UR3	C2-N3	5.32	1.49	1.39
55	X	56	OMC	C4-N3	5.31	1.45	1.34
1	A	967	5MC	C6-N1	5.30	1.47	1.38
55	X	48	5MC	C6-N1	5.21	1.46	1.38
22	a	2503	2MA	C5-C6	5.21	1.55	1.41
1	A	1407	5MC	C6-N1	5.18	1.46	1.38
53	V	55	PSU	C6-N1	5.16	1.44	1.36
22	a	2504	PSU	C6-N1	5.13	1.44	1.36
1	A	966	2MG	C2-N1	5.11	1.44	1.36
22	a	1962	5MC	C6-N1	5.10	1.46	1.38
22	a	746	PSU	C6-N1	5.09	1.44	1.36
22	a	2503	2MA	C6-N6	-5.04	1.21	1.34
1	A	1207	2MG	C2-N1	5.02	1.44	1.36
55	X	55	PSU	C6-N1	5.01	1.44	1.36
55	X	56	OMC	C4-N4	5.01	1.45	1.33
53	V	13	PSU	C6-N1	5.00	1.44	1.36
1	A	516	PSU	C6-N1	5.00	1.44	1.36
22	a	2251	OMG	C2-N2	4.94	1.45	1.34
22	a	2605	PSU	C6-N1	4.93	1.44	1.36
22	a	2604	PSU	C6-N1	4.92	1.44	1.36
22	a	1917	PSU	C6-N1	4.86	1.44	1.36
55	X	57	7S3	C2-N1	4.82	1.44	1.35
55	X	15	7SN	C2-N2	4.81	1.45	1.34
22	a	1835	2MG	C2-N1	4.81	1.44	1.36
55	X	37	T6A	C10-N11	4.80	1.45	1.35
22	a	1911	PSU	C6-N1	4.80	1.44	1.36
22	a	2457	PSU	C6-N1	4.80	1.44	1.36
53	V	38	2MA	C6-N6	-4.78	1.22	1.34
22	a	2498	OMC	C4-N3	4.78	1.44	1.34
1	A	1516	2MG	C2-N1	4.77	1.44	1.36
22	a	2580	PSU	C6-N1	4.74	1.44	1.36
55	X	56	OMC	C2-N1	4.73	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	2498	OMC	C4-N4	4.70	1.45	1.33
53	V	35	U8U	C6-N1	4.69	1.46	1.38
22	a	2445	2MG	C2-N1	4.69	1.44	1.36
22	a	2498	OMC	C2-N1	4.60	1.50	1.40
22	a	955	PSU	C6-N1	4.52	1.43	1.36
1	A	1402	4OC	C4-N4	4.49	1.45	1.35
22	a	2069	G7M	C5-N7	-4.47	1.34	1.39
22	a	1915	3TD	C2-N3	4.44	1.48	1.38
55	X	15	7SN	C5-C6	4.34	1.53	1.44
55	X	15	7SN	CA7-N71	4.29	1.45	1.28
1	A	1407	5MC	C2-N1	4.24	1.49	1.40
55	X	15	7SN	C7-CA7	4.22	1.53	1.46
1	A	1402	4OC	C2-N1	4.15	1.49	1.40
1	A	527	G7M	C5-N7	-4.15	1.34	1.39
55	X	48	5MC	C4-N4	4.12	1.44	1.34
55	X	48	5MC	C2-N1	4.11	1.48	1.40
1	A	967	5MC	C4-N4	4.10	1.44	1.34
1	A	967	5MC	C2-N1	3.99	1.48	1.40
1	A	1407	5MC	C4-N4	3.96	1.44	1.34
1	A	1402	4OC	C5-C4	3.91	1.49	1.40
22	a	1962	5MC	C4-N4	3.87	1.44	1.34
33	l	82	MS6	CB-CA	3.85	1.58	1.53
1	A	527	G7M	C5-C6	3.85	1.54	1.43
53	V	35	U8U	C4-N3	3.84	1.46	1.38
22	a	1962	5MC	C2-N1	3.83	1.48	1.40
22	a	2069	G7M	C5-C6	3.79	1.54	1.43
22	a	2503	2MA	C6-N1	3.76	1.40	1.35
55	X	15	7SN	CA7-N72	-3.60	1.25	1.33
53	V	35	U8U	C2-S2	-3.58	1.62	1.67
55	X	57	7S3	C5-C6	3.57	1.54	1.45
55	X	54	B8H	C6-N1	3.53	1.45	1.36
53	V	38	2MA	C6-N1	3.51	1.40	1.35
53	V	13	PSU	C4-N3	3.40	1.45	1.38
55	X	55	PSU	C4-N3	3.38	1.45	1.38
12	L	89	D2T	CB-SB	3.29	1.85	1.82
53	V	55	PSU	C4-N3	3.29	1.45	1.38
1	A	516	PSU	C4-N3	3.28	1.44	1.38
22	a	1911	PSU	C4-N3	3.26	1.44	1.38
22	a	746	PSU	C4-N3	3.25	1.44	1.38
1	A	1402	4OC	O2-C2	-3.23	1.17	1.23
22	a	2457	PSU	C4-N3	3.22	1.44	1.38
55	X	54	B8H	C1'-C5	3.20	1.57	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	2445	2MG	C5-N7	-3.18	1.32	1.39
22	a	745	1MG	C5-C6	3.17	1.53	1.45
1	A	1402	4OC	C6-N1	3.17	1.45	1.38
22	a	2604	PSU	C4-N3	3.16	1.44	1.38
55	X	56	OMC	C6-N1	3.16	1.45	1.38
22	a	2251	OMG	C5-N7	-3.15	1.33	1.39
53	V	38	2MA	C5-N7	-3.15	1.33	1.39
22	a	955	PSU	C4-N3	3.14	1.44	1.38
22	a	2503	2MA	C5-N7	-3.13	1.33	1.39
33	l	81	4D4	CZ-NE	3.13	1.39	1.33
22	a	2498	OMC	C6-N1	3.12	1.45	1.38
12	L	89	D2T	CB-CG	3.11	1.57	1.52
1	A	1207	2MG	C5-N7	-3.10	1.33	1.39
1	A	1519	MA6	C6-N6	3.10	1.46	1.36
22	a	1835	2MG	C5-N7	-3.10	1.33	1.39
22	a	1917	PSU	C4-N3	3.09	1.44	1.38
22	a	2504	PSU	C4-N3	3.08	1.44	1.38
22	a	2251	OMG	O6-C6	-3.08	1.17	1.23
1	A	1518	MA6	C6-N6	3.08	1.45	1.36
22	a	2552	OMU	C4-N3	3.07	1.44	1.38
1	A	1498	UR3	C6-N1	3.06	1.45	1.38
1	A	966	2MG	C5-N7	-3.03	1.33	1.39
22	a	745	1MG	C5-N7	-3.00	1.33	1.39
22	a	2498	OMC	O2-C2	-3.00	1.18	1.23
1	A	527	G7M	C6-N1	2.97	1.44	1.38
22	a	2580	PSU	C4-N3	2.97	1.44	1.38
22	a	2605	PSU	C4-N3	2.97	1.44	1.38
53	V	35	U8U	C4-C5	2.96	1.51	1.45
1	A	1518	MA6	C5-C4	-2.95	1.33	1.39
22	a	2449	H2U	C2-N3	-2.94	1.32	1.38
1	A	1519	MA6	C5-C4	-2.93	1.33	1.39
55	X	56	OMC	O2-C2	-2.93	1.18	1.23
1	A	527	G7M	C2-N1	2.92	1.44	1.37
22	a	1915	3TD	C4-N3	2.91	1.46	1.40
1	A	1516	2MG	C5-N7	-2.88	1.33	1.39
22	a	2552	OMU	O4-C4	-2.88	1.18	1.24
55	X	57	7S3	C6-N1	2.87	1.43	1.39
22	a	2552	OMU	O2-C2	-2.87	1.17	1.23
22	a	2449	H2U	C4-N3	-2.86	1.32	1.37
22	a	2251	OMG	C5-C6	2.80	1.54	1.44
55	X	54	B8H	O2-C2	-2.79	1.18	1.23
1	A	1407	5MC	O2-C2	-2.75	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	2069	G7M	C2-N1	2.73	1.44	1.37
22	a	2069	G7M	O6-C6	-2.71	1.18	1.23
22	a	2030	6MZ	C5-C4	-2.71	1.34	1.39
22	a	1835	2MG	O6-C6	-2.69	1.18	1.23
1	A	1516	2MG	C5-C6	2.69	1.54	1.44
22	a	2069	G7M	C6-N1	2.68	1.43	1.38
1	A	966	2MG	C5-C6	2.68	1.54	1.44
1	A	1207	2MG	C5-C6	2.67	1.54	1.44
22	a	2445	2MG	C5-C6	2.65	1.54	1.44
22	a	2445	2MG	O6-C6	-2.65	1.18	1.23
55	X	15	7SN	C2-N1	2.65	1.44	1.37
1	A	1518	MA6	C5-N7	-2.64	1.34	1.39
55	X	57	7S3	C5-N7	-2.58	1.34	1.39
1	A	1519	MA6	C5-N7	-2.57	1.34	1.39
22	a	1835	2MG	C5-C6	2.55	1.53	1.44
22	a	1939	5MU	O2-C2	-2.52	1.18	1.23
1	A	966	2MG	C6-N1	2.52	1.43	1.38
1	A	1516	2MG	O6-C6	-2.51	1.18	1.23
55	X	15	7SN	O6-C6	-2.51	1.18	1.23
22	a	1962	5MC	O2-C2	-2.50	1.19	1.23
55	X	54	B8H	O4-C4	-2.49	1.18	1.23
55	X	15	7SN	C7-C5	-2.49	1.37	1.44
1	A	1498	UR3	O4-C4	-2.48	1.18	1.23
22	a	1618	6MZ	C5-C4	-2.47	1.34	1.39
22	a	2251	OMG	C2-N1	2.47	1.43	1.37
55	X	56	OMC	C5-C4	2.46	1.48	1.42
22	a	2498	OMC	C5-C4	2.46	1.48	1.42
22	a	745	1MG	C4-N9	-2.45	1.31	1.38
1	A	527	G7M	O6-C6	-2.45	1.18	1.23
1	A	967	5MC	O2-C2	-2.44	1.19	1.23
22	a	745	1MG	O6-C6	-2.43	1.18	1.23
22	a	2604	PSU	O4-C4	-2.42	1.19	1.23
1	A	1207	2MG	C6-N1	2.41	1.43	1.38
22	a	746	PSU	O4-C4	-2.41	1.19	1.23
22	a	1618	6MZ	C5-N7	-2.39	1.34	1.39
22	a	2503	2MA	CM2-C2	2.37	1.56	1.49
22	a	2251	OMG	C6-N1	2.36	1.43	1.38
22	a	2605	PSU	O4-C4	-2.35	1.19	1.23
22	a	2457	PSU	O4-C4	-2.35	1.19	1.23
1	A	1519	MA6	C8-N9	-2.33	1.33	1.37
55	X	37	T6A	C5-C4	-2.32	1.34	1.39
1	A	1207	2MG	O6-C6	-2.32	1.19	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	966	2MG	O6-C6	-2.31	1.19	1.23
53	V	38	2MA	CM2-C2	2.31	1.56	1.49
55	X	48	5MC	CM5-C5	2.29	1.56	1.50
55	X	26	M2G	C2-N2	2.28	1.39	1.35
22	a	746	PSU	O4'-C1'	-2.27	1.40	1.43
22	a	747	5MU	O2-C2	-2.27	1.18	1.23
1	A	516	PSU	O4-C4	-2.27	1.19	1.23
22	a	2580	PSU	O4'-C1'	-2.26	1.40	1.43
1	A	967	5MC	CM5-C5	2.25	1.56	1.50
22	a	1917	PSU	O4-C4	-2.25	1.19	1.23
1	A	1498	UR3	O2-C2	-2.25	1.18	1.22
1	A	1498	UR3	C5-C4	2.25	1.49	1.43
55	X	15	7SN	C6-N1	2.24	1.43	1.38
22	a	2580	PSU	O4-C4	-2.24	1.19	1.23
1	A	1516	2MG	C6-N1	2.23	1.43	1.38
22	a	955	PSU	O4-C4	-2.23	1.19	1.23
22	a	2504	PSU	O4-C4	-2.22	1.19	1.23
22	a	2030	6MZ	C5-N7	-2.20	1.34	1.39
22	a	1962	5MC	CM5-C5	2.20	1.56	1.50
1	A	1518	MA6	C8-N9	-2.19	1.33	1.37
22	a	1911	PSU	O4-C4	-2.17	1.19	1.23
55	X	37	T6A	C6-N1	-2.15	1.31	1.35
22	a	2445	2MG	C6-N1	2.14	1.42	1.38
55	X	48	5MC	O2-C2	-2.13	1.19	1.23
22	a	2552	OMU	C5-C4	2.13	1.48	1.43
1	A	1407	5MC	CM5-C5	2.13	1.56	1.50
22	a	1835	2MG	C6-N1	2.13	1.42	1.38
22	a	1915	3TD	O4-C4	-2.12	1.18	1.23
55	X	55	PSU	O4-C4	-2.12	1.19	1.23
22	a	2030	6MZ	C8-N9	-2.11	1.33	1.37
53	V	55	PSU	O4-C4	-2.10	1.19	1.23
22	a	955	PSU	C4-C5	2.09	1.50	1.44
55	X	37	T6A	O10-C10	-2.09	1.19	1.23
33	l	81	4D4	OB-CB	-2.08	1.39	1.43
55	X	37	T6A	ODB-C13	-2.06	1.23	1.30
1	A	1498	UR3	C4-N3	2.06	1.45	1.40
1	A	1516	2MG	C4-N9	-2.05	1.32	1.38
22	a	1939	5MU	O4-C4	-2.04	1.19	1.23
1	A	527	G7M	C8-N7	2.03	1.36	1.33
55	X	57	7S3	O6-C6	-2.01	1.18	1.23
22	a	2445	2MG	C4-N9	-2.01	1.32	1.38
53	V	13	PSU	O4-C4	-2.01	1.19	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	a	1618	6MZ	C8-N9	-2.00	1.33	1.37

All (296) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	a	747	5MU	C5-C4-N3	12.45	125.94	115.31
53	V	54	5MU	C5-C4-N3	12.00	125.56	115.31
22	a	1939	5MU	C5-C4-N3	11.98	125.54	115.31
22	a	1939	5MU	C5-C6-N1	-10.12	112.93	123.34
22	a	747	5MU	C5-C6-N1	-9.94	113.11	123.34
53	V	54	5MU	C5-C6-N1	-9.24	113.83	123.34
22	a	2503	2MA	C5-C4-N3	-8.31	117.84	127.19
53	V	38	2MA	C4-N9-C1'	-7.96	107.64	126.59
53	V	38	2MA	C1'-N9-C8	7.71	144.55	127.14
53	V	38	2MA	C5-C4-N3	-7.21	119.08	127.19
1	A	966	2MG	C2-N3-C4	7.04	120.77	112.04
22	a	2445	2MG	C2-N3-C4	6.99	120.70	112.04
22	a	1835	2MG	C2-N3-C4	6.63	120.26	112.04
22	a	2503	2MA	C4-N9-C1'	-6.48	111.15	126.59
1	A	1207	2MG	C2-N3-C4	6.46	120.05	112.04
1	A	1516	2MG	C2-N3-C4	6.38	119.94	112.04
22	a	2503	2MA	C1'-N9-C8	6.35	141.46	127.14
1	A	1518	MA6	N1-C2-N3	-6.07	119.10	128.60
55	X	15	7SN	C5-C4-N3	-5.95	120.56	127.52
22	a	2030	6MZ	N1-C2-N3	-5.89	119.38	128.60
22	a	2503	2MA	N3-C4-N9	5.78	135.01	126.99
1	A	1519	MA6	N1-C2-N3	-5.76	119.58	128.60
55	X	37	T6A	N3-C2-N1	-5.76	119.59	128.60
22	a	747	5MU	O4-C4-C5	-5.53	118.49	124.90
1	A	966	2MG	C5-C4-N3	-5.51	119.52	128.46
22	a	1618	6MZ	C5-C4-N3	-5.50	119.58	126.75
22	a	2030	6MZ	C9-N6-C6	-5.49	118.14	122.87
22	a	2251	OMG	C5-C4-N3	-5.49	119.56	128.46
55	X	37	T6A	N6-C10-N11	5.31	121.18	113.76
55	X	37	T6A	C5-C4-N3	-5.31	119.83	126.75
22	a	1835	2MG	C5-C4-N3	-5.28	119.90	128.46
22	a	1618	6MZ	N1-C2-N3	-5.26	120.38	128.60
22	a	2552	OMU	C4-N3-C2	-5.26	119.65	126.58
55	X	57	7S3	C5-C4-N3	-5.13	119.60	127.26
22	a	2445	2MG	C5-C4-N3	-5.12	120.15	128.46
22	a	1939	5MU	C4-N3-C2	-4.99	120.89	127.35
22	a	1939	5MU	N3-C2-N1	4.98	121.50	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1498	UR3	C4-N3-C2	-4.97	119.88	124.56
22	a	747	5MU	C4-N3-C2	-4.90	121.01	127.35
1	A	1207	2MG	C5-C4-N3	-4.88	120.54	128.46
53	V	54	5MU	O4-C4-C5	-4.86	119.27	124.90
22	a	955	PSU	C4-N3-C2	-4.86	119.34	126.34
22	a	1915	3TD	N1-C2-N3	4.84	119.96	116.14
22	a	2030	6MZ	C5-C4-N3	-4.82	120.47	126.75
22	a	2457	PSU	C4-N3-C2	-4.79	119.44	126.34
1	A	1518	MA6	N9-C8-N7	-4.77	107.40	113.91
1	A	1518	MA6	C2-N1-C6	4.75	122.97	111.75
55	X	57	7S3	N1-C2-N3	-4.75	120.36	126.00
22	a	2604	PSU	C4-N3-C2	-4.69	119.58	126.34
1	A	1519	MA6	N9-C8-N7	-4.69	107.50	113.91
22	a	2605	PSU	C4-N3-C2	-4.69	119.59	126.34
53	V	55	PSU	C4-N3-C2	-4.68	119.59	126.34
22	a	2030	6MZ	N9-C8-N7	-4.68	107.51	113.91
1	A	1516	2MG	C5-C4-N3	-4.68	120.87	128.46
22	a	746	PSU	C4-N3-C2	-4.65	119.64	126.34
22	a	1917	PSU	C4-N3-C2	-4.65	119.64	126.34
22	a	955	PSU	N1-C2-N3	4.62	120.36	115.13
53	V	13	PSU	C4-N3-C2	-4.58	119.73	126.34
1	A	516	PSU	C4-N3-C2	-4.58	119.74	126.34
22	a	2457	PSU	N1-C2-N3	4.58	120.31	115.13
53	V	54	5MU	C5M-C5-C4	4.57	123.80	118.77
1	A	1519	MA6	C2-N1-C6	4.53	122.46	111.75
53	V	38	2MA	N3-C4-N9	4.53	133.27	126.99
22	a	2504	PSU	C4-N3-C2	-4.52	119.83	126.34
1	A	1519	MA6	C5-C4-N3	-4.52	120.86	126.75
22	a	1911	PSU	C4-N3-C2	-4.51	119.84	126.34
53	V	54	5MU	C4-N3-C2	-4.51	121.52	127.35
53	V	55	PSU	N1-C2-N3	4.46	120.19	115.13
22	a	1618	6MZ	C4-C5-C6	4.44	120.25	116.81
55	X	15	7SN	C1'-N9-C4	-4.43	113.35	126.50
22	a	2069	G7M	C2-N3-C4	4.43	120.19	112.30
1	A	966	2MG	C2-N1-C6	-4.42	119.39	124.48
22	a	745	1MG	C5-C4-N3	-4.42	121.30	128.46
55	X	54	B8H	C4-N3-C2	-4.41	121.64	127.35
22	a	2251	OMG	C2-N3-C4	4.41	120.16	112.30
55	X	55	PSU	C4-N3-C2	-4.41	119.98	126.34
55	X	15	7SN	C2-N3-C4	4.41	120.15	112.30
55	X	37	T6A	N9-C8-N7	-4.39	107.91	113.91
22	a	2504	PSU	N1-C2-N3	4.37	120.08	115.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	a	2580	PSU	N1-C2-N3	4.36	120.07	115.13
22	a	2605	PSU	N1-C2-N3	4.36	120.07	115.13
53	V	54	5MU	N3-C2-N1	4.33	120.63	114.89
1	A	1518	MA6	C5-C4-N3	-4.32	121.11	126.75
22	a	1939	5MU	O4-C4-C5	-4.29	119.93	124.90
22	a	1618	6MZ	N9-C8-N7	-4.27	108.07	113.91
1	A	1207	2MG	C2-N1-C6	-4.26	119.58	124.48
1	A	1519	MA6	C4-C5-C6	4.25	120.63	115.88
22	a	1835	2MG	C2-N1-C6	-4.21	119.63	124.48
1	A	527	G7M	C2-N3-C4	4.21	119.80	112.30
22	a	1917	PSU	N1-C2-N3	4.20	119.89	115.13
53	V	54	5MU	C5M-C5-C6	-4.19	117.25	122.85
22	a	2580	PSU	C4-N3-C2	-4.15	120.36	126.34
22	a	747	5MU	N3-C2-N1	4.14	120.39	114.89
22	a	2604	PSU	N1-C2-N3	4.13	119.81	115.13
55	X	55	PSU	N1-C2-N3	4.06	119.72	115.13
55	X	37	T6A	C4-C5-C6	4.02	119.92	116.81
53	V	38	2MA	N9-C8-N7	-3.99	108.46	113.91
1	A	1518	MA6	C4-C5-C6	3.97	120.33	115.88
53	V	13	PSU	N1-C2-N3	3.97	119.63	115.13
22	a	1911	PSU	N1-C2-N3	3.96	119.61	115.13
22	a	2445	2MG	C2-N1-C6	-3.95	119.94	124.48
1	A	516	PSU	N1-C2-N3	3.93	119.58	115.13
22	a	1915	3TD	C4-N3-C2	-3.92	120.35	124.61
22	a	2069	G7M	C5-C6-N1	3.91	119.96	111.79
22	a	747	5MU	C5M-C5-C4	3.90	123.06	118.77
55	X	57	7S3	C2-N3-C4	3.89	120.05	112.41
33	l	81	4D4	NE-CZ-NH2	3.88	127.52	120.70
22	a	747	5MU	C5M-C5-C6	-3.86	117.69	122.85
1	A	527	G7M	C5-C6-N1	3.85	119.84	111.79
22	a	2251	OMG	N9-C4-N3	3.83	133.63	125.94
22	a	2552	OMU	N3-C2-N1	3.83	119.97	114.89
22	a	746	PSU	N1-C2-N3	3.82	119.46	115.13
1	A	527	G7M	C5-C4-N3	-3.80	120.87	128.15
1	A	1516	2MG	C2-N1-C6	-3.78	120.13	124.48
1	A	1519	MA6	C5-N7-C8	3.77	108.86	103.51
1	A	1518	MA6	C5-N7-C8	3.76	108.86	103.51
55	X	48	5MC	C5-C6-N1	-3.76	119.47	123.34
22	a	2069	G7M	C5-C4-N3	-3.72	121.01	128.15
22	a	1939	5MU	C5M-C5-C6	-3.72	117.88	122.85
1	A	966	2MG	N9-C4-N3	3.68	133.33	125.94
22	a	745	1MG	C2-N3-C4	3.66	120.20	111.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	a	2030	6MZ	C2-N3-C4	3.65	120.37	111.75
22	a	2580	PSU	C6-N1-C2	-3.59	119.01	122.68
55	X	37	T6A	C2-N3-C4	3.59	120.24	111.75
1	A	1407	5MC	C5-C6-N1	-3.59	119.65	123.34
22	a	1618	6MZ	N3-C4-N9	3.58	132.99	127.08
22	a	2503	2MA	N9-C8-N7	-3.54	109.08	113.91
1	A	967	5MC	C5-C6-N1	-3.53	119.71	123.34
22	a	1618	6MZ	C2-N3-C4	3.51	120.04	111.75
22	a	1939	5MU	C5M-C5-C4	3.46	122.58	118.77
1	A	1516	2MG	N9-C8-N7	-3.45	106.89	113.39
22	a	1835	2MG	N9-C4-N3	3.44	132.85	125.94
22	a	2445	2MG	N9-C8-N7	-3.40	106.98	113.39
55	X	15	7SN	C7-CA7-N72	3.40	121.34	117.69
22	a	2030	6MZ	C5-N7-C8	3.40	108.33	103.51
22	a	2552	OMU	C5-C4-N3	3.38	119.90	114.84
53	V	55	PSU	C6-N1-C2	-3.36	119.24	122.68
1	A	527	G7M	O6-C6-C5	-3.36	120.49	128.06
22	a	2251	OMG	N9-C8-N7	-3.29	107.19	113.39
55	X	57	7S3	N9-C8-N7	-3.27	107.23	113.39
1	A	1207	2MG	N9-C8-N7	-3.25	107.26	113.39
22	a	2251	OMG	C2-N1-C6	-3.25	119.18	125.10
55	X	37	T6A	N3-C4-N9	3.24	132.43	127.08
55	X	37	T6A	C5-N7-C8	3.24	108.11	103.51
22	a	2069	G7M	O6-C6-C5	-3.23	120.76	128.06
1	A	1518	MA6	C2-N3-C4	3.23	119.37	111.75
22	a	2503	2MA	N3-C2-N1	-3.23	119.79	125.72
22	a	2445	2MG	N9-C4-N3	3.22	132.41	125.94
1	A	1519	MA6	C2-N3-C4	3.21	119.34	111.75
22	a	745	1MG	N9-C8-N7	-3.21	107.35	113.39
55	X	57	7S3	C1'-N9-C4	-3.19	117.02	126.50
22	a	2504	PSU	C6-N1-C2	-3.18	119.43	122.68
53	V	38	2MA	C5-N7-C8	3.16	108.00	103.51
22	a	1962	5MC	C5-C6-N1	-3.16	120.09	123.34
22	a	1618	6MZ	C5-N7-C8	3.15	107.98	103.51
1	A	966	2MG	N9-C8-N7	-3.12	107.52	113.39
12	L	89	D2T	OD2-CG-CB	3.10	119.84	113.15
22	a	2580	PSU	C6-C5-C4	3.07	120.34	118.20
53	V	35	U8U	C5-C4-N3	3.06	119.28	114.97
53	V	13	PSU	C6-C5-C4	3.06	120.34	118.20
22	a	1917	PSU	C6-C5-C4	3.04	120.32	118.20
22	a	2457	PSU	C6-N1-C2	-3.04	119.58	122.68
1	A	1519	MA6	C4-C5-N7	-3.02	106.95	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	a	955	PSU	O2-C2-N1	-2.99	119.50	122.79
33	l	82	MS6	CE-SD-CG	2.98	110.64	100.40
1	A	1207	2MG	N9-C4-N3	2.97	131.91	125.94
1	A	527	G7M	N9-C4-N3	2.97	131.90	125.94
22	a	1835	2MG	N9-C8-N7	-2.96	107.81	113.39
22	a	2457	PSU	C6-C5-C4	2.96	120.27	118.20
55	X	55	PSU	C6-N1-C2	-2.96	119.66	122.68
22	a	2605	PSU	C6-N1-C2	-2.95	119.67	122.68
55	X	37	T6A	O10-C10-N6	-2.95	118.64	123.62
22	a	1939	5MU	O4-C4-N3	-2.91	114.53	120.12
53	V	38	2MA	N3-C2-N1	-2.90	120.38	125.72
22	a	2445	2MG	N1-C2-N3	-2.90	119.46	123.95
22	a	2503	2MA	C6-C5-C4	2.90	121.08	117.18
1	A	1518	MA6	C4-C5-N7	-2.90	107.09	110.62
1	A	1516	2MG	N9-C4-N3	2.89	131.74	125.94
22	a	745	1MG	C5-C6-N1	2.89	120.49	114.91
22	a	955	PSU	C6-N1-C2	-2.84	119.78	122.68
55	X	57	7S3	N9-C4-N3	2.83	133.55	126.94
22	a	1911	PSU	C6-C5-C4	2.81	120.17	118.20
22	a	1939	5MU	O2-C2-N1	-2.81	119.05	122.79
22	a	2030	6MZ	C4-C5-C6	2.81	118.98	116.81
22	a	2251	OMG	C5-C6-N1	2.81	120.32	113.19
1	A	527	G7M	C2-N1-C6	-2.80	119.99	125.10
1	A	966	2MG	C5-C6-N1	2.78	120.26	113.19
22	a	2504	PSU	C6-C5-C4	2.77	120.14	118.20
1	A	1516	2MG	C5-C6-N1	2.77	120.23	113.19
53	V	38	2MA	C6-C5-C4	2.77	120.91	117.18
22	a	1835	2MG	C5-C6-N1	2.77	120.22	113.19
22	a	2503	2MA	C5-N7-C8	2.77	107.44	103.51
22	a	2069	G7M	N9-C4-N3	2.76	131.49	125.94
22	a	955	PSU	C6-C5-C4	2.76	120.13	118.20
1	A	1516	2MG	N1-C2-N3	-2.75	119.70	123.95
22	a	2445	2MG	C5-C6-N1	2.73	120.13	113.19
1	A	527	G7M	CN7-N7-C5	2.72	130.14	126.77
22	a	1917	PSU	C6-N1-C2	-2.71	119.91	122.68
22	a	2069	G7M	C2-N1-C6	-2.70	120.18	125.10
22	a	1618	6MZ	C9-N6-C6	-2.69	120.56	122.87
1	A	1207	2MG	C5-C6-N1	2.67	119.97	113.19
22	a	2605	PSU	C6-C5-C4	2.67	120.06	118.20
12	L	89	D2T	OD1-CG-CB	-2.65	116.88	122.44
55	X	55	PSU	O2-C2-N1	-2.65	119.88	122.79
22	a	2604	PSU	C6-C5-C4	2.64	120.04	118.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	966	2MG	N1-C2-N3	-2.63	119.88	123.95
22	a	2580	PSU	O4'-C1'-C2'	2.63	108.85	105.14
55	X	15	7SN	C2-N1-C6	-2.60	120.35	125.10
22	a	2580	PSU	O2-C2-N1	-2.59	119.94	122.79
1	A	1516	2MG	C1'-N9-C4	-2.59	118.82	126.50
53	V	54	5MU	O4-C4-N3	-2.58	115.18	120.12
22	a	2030	6MZ	N3-C4-N9	2.57	131.32	127.08
22	a	2604	PSU	C6-N1-C2	-2.57	120.06	122.68
55	X	15	7SN	N9-C4-N3	2.57	131.09	125.94
1	A	966	2MG	O6-C6-C5	-2.56	119.81	126.60
1	A	516	PSU	C6-N1-C2	-2.56	120.07	122.68
33	l	81	4D4	O-C-CA	-2.55	118.09	124.78
22	a	1835	2MG	N1-C2-N3	-2.55	120.01	123.95
55	X	57	7S3	C1'-N9-C8	2.54	133.92	126.70
22	a	2251	OMG	O6-C6-C5	-2.52	119.91	126.60
22	a	2457	PSU	O2-C2-N1	-2.51	120.02	122.79
33	l	81	4D4	CB-CA-C	-2.51	107.77	111.77
1	A	1207	2MG	O6-C6-C5	-2.51	119.95	126.60
22	a	2030	6MZ	C4-C5-N7	-2.50	107.58	110.62
22	a	2552	OMU	O4-C4-C5	-2.49	120.79	125.16
22	a	1911	PSU	C6-N1-C2	-2.48	120.15	122.68
1	A	1402	4OC	C6-C5-C4	2.47	119.98	116.96
53	V	13	PSU	C6-N1-C2	-2.47	120.16	122.68
1	A	1518	MA6	N1-C6-N6	-2.46	114.39	117.08
22	a	1835	2MG	O6-C6-C5	-2.46	120.06	126.60
22	a	1911	PSU	O2-C2-N1	-2.46	120.09	122.79
1	A	516	PSU	C6-C5-C4	2.46	119.92	118.20
1	A	1207	2MG	N1-C2-N3	-2.44	120.18	123.95
53	V	55	PSU	C6-C5-C4	2.44	119.90	118.20
22	a	2030	6MZ	C4-N9-C8	2.44	108.37	105.73
22	a	746	PSU	C6-N1-C2	-2.43	120.20	122.68
22	a	2445	2MG	C8-N7-C5	2.42	108.62	104.24
53	V	55	PSU	O2-C2-N1	-2.41	120.13	122.79
55	X	57	7S3	C8-N7-C5	2.40	108.58	104.24
22	a	747	5MU	O2-C2-N1	-2.40	119.60	122.79
22	a	2552	OMU	O2-C2-N1	-2.38	119.62	122.79
22	a	747	5MU	O4-C4-N3	-2.37	115.58	120.12
1	A	1407	5MC	CM5-C5-C6	-2.36	119.69	122.85
53	V	13	PSU	O2-C2-N1	-2.36	120.19	122.79
53	V	35	U8U	C1'-N1-C6	-2.36	117.20	121.12
1	A	1518	MA6	C4-N9-C8	2.35	108.28	105.73
1	A	1207	2MG	C8-N7-C5	2.35	108.50	104.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1516	2MG	C8-N7-C5	2.35	108.50	104.24
25	d	150	MEQ	OE1-CD-CG	2.35	126.31	122.02
1	A	1519	MA6	C5-C4-N9	2.34	108.50	105.78
22	a	2605	PSU	O2-C2-N1	-2.34	120.22	122.79
53	V	35	U8U	O4-C4-C5	-2.32	121.51	124.96
22	a	2251	OMG	C8-N7-C5	2.31	108.41	104.24
12	L	89	D2T	O-C-CA	-2.29	118.77	124.78
55	X	55	PSU	C6-C5-C4	2.29	119.80	118.20
1	A	1498	UR3	C1'-N1-C2	2.26	120.81	116.99
1	A	1519	MA6	C4-N9-C8	2.26	108.18	105.73
1	A	966	2MG	C8-N7-C5	2.26	108.33	104.24
22	a	1618	6MZ	C4-N9-C8	2.26	108.17	105.73
22	a	2069	G7M	CN7-N7-C5	2.24	129.55	126.77
1	A	1518	MA6	C5-C4-N9	2.23	108.38	105.78
1	A	1498	UR3	C6-N1-C2	-2.23	119.79	121.79
22	a	745	1MG	C8-N7-C5	2.22	108.26	104.24
1	A	516	PSU	O4'-C1'-C2'	2.22	108.27	105.14
22	a	2445	2MG	O6-C6-C5	-2.21	120.73	126.60
1	A	516	PSU	O2-C2-N1	-2.20	120.36	122.79
22	a	1917	PSU	O2-C2-N1	-2.20	120.37	122.79
22	a	745	1MG	N9-C4-N3	2.19	130.34	125.94
22	a	1962	5MC	CM5-C5-C6	-2.18	119.94	122.85
22	a	2604	PSU	O2-C2-N1	-2.18	120.40	122.79
55	X	37	T6A	C4-N9-C8	2.16	108.07	105.73
22	a	2069	G7M	N9-C8-N7	-2.16	106.88	112.21
1	A	1519	MA6	N3-C4-N9	2.16	130.63	127.08
1	A	1516	2MG	O6-C6-C5	-2.15	120.89	126.60
22	a	2251	OMG	C1'-N9-C4	-2.14	120.15	126.50
55	X	26	M2G	O6-C6-N1	-2.14	116.02	120.12
22	a	2503	2MA	N6-C6-N1	2.13	120.05	117.03
25	d	150	MEQ	CG-CD-NE2	-2.13	113.33	116.29
55	X	26	M2G	CM2-N2-CM1	2.12	122.33	115.77
11	K	119	IAS	OXT-C-O	-2.11	119.29	124.09
1	A	1402	4OC	O2-C2-N3	-2.11	118.89	122.33
22	a	2069	G7M	N1-C2-N3	-2.11	119.39	123.32
11	K	119	IAS	OXT-C-CA	2.10	120.54	113.38
33	l	81	4D4	NH1-CZ-NE	-2.10	114.34	119.19
53	V	38	2MA	N6-C6-N1	2.10	120.00	117.03
22	a	747	5MU	C1'-N1-C2	2.10	121.36	117.57
22	a	2030	6MZ	C5-C4-N9	2.09	108.21	105.78
1	A	1518	MA6	N3-C4-N9	2.08	130.51	127.08
53	V	38	2MA	C4-C5-N7	-2.07	108.09	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	a	1835	2MG	C8-N7-C5	2.06	107.97	104.24
1	A	527	G7M	CN7-N7-C8	-2.05	121.67	124.84
22	a	2457	PSU	O4'-C1'-C2'	2.04	108.02	105.14
55	X	57	7S3	C2-N1-C6	-2.03	119.42	122.11
55	X	37	T6A	C4-C5-N7	-2.01	108.17	110.62
22	a	745	1MG	C4-C5-N7	-2.01	107.54	110.72
1	A	1407	5MC	O2-C2-N3	-2.00	119.08	122.33

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	L	89	D2T	SB-CB-CG-OD2
25	d	150	MEQ	C-CA-CB-CG
22	a	2251	OMG	C1'-C2'-O2'-CM2
55	X	26	M2G	O4'-C4'-C5'-O5'
55	X	26	M2G	C3'-C4'-C5'-O5'
55	X	48	5MC	O4'-C4'-C5'-O5'
55	X	48	5MC	C3'-C4'-C5'-O5'
22	a	1915	3TD	C3'-C4'-C5'-O5'
22	a	1915	3TD	O4'-C4'-C5'-O5'
22	a	2030	6MZ	O4'-C4'-C5'-O5'
22	a	2504	PSU	C3'-C4'-C5'-O5'
22	a	2504	PSU	O4'-C4'-C5'-O5'
53	V	13	PSU	C3'-C4'-C5'-O5'
53	V	54	5MU	C3'-C4'-C5'-O5'
53	V	54	5MU	O4'-C4'-C5'-O5'
1	A	1519	MA6	O4'-C4'-C5'-O5'
55	X	57	7S3	C3'-C4'-C5'-O5'
55	X	57	7S3	O4'-C4'-C5'-O5'
25	d	150	MEQ	OE1-CD-CG-CB
25	d	150	MEQ	NE2-CD-CG-CB
55	X	15	7SN	C3'-C4'-C5'-O5'
55	X	37	T6A	N11-C12-C14-O14
22	a	2030	6MZ	C3'-C4'-C5'-O5'
55	X	37	T6A	N11-C12-C13-ODB
1	A	966	2MG	C3'-C4'-C5'-O5'
53	V	13	PSU	O4'-C4'-C5'-O5'
55	X	15	7SN	O4'-C4'-C5'-O5'
1	A	1519	MA6	C3'-C4'-C5'-O5'
55	X	37	T6A	N11-C12-C14-C15
1	A	1402	4OC	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
22	a	2445	2MG	C3'-C4'-C5'-O5'
55	X	37	T6A	C13-C12-C14-O14
55	X	37	T6A	C13-C12-C14-C15
1	A	966	2MG	O4'-C4'-C5'-O5'
22	a	2552	OMU	C3'-C2'-O2'-CM2
25	d	150	MEQ	N-CA-CB-CG
1	A	527	G7M	C3'-C4'-C5'-O5'
1	A	527	G7M	C4'-C5'-O5'-P
55	X	37	T6A	N11-C12-C13-ODA
53	V	55	PSU	O4'-C1'-C5-C4
55	X	55	PSU	O4'-C1'-C5-C4
1	A	1402	4OC	C3'-C4'-C5'-O5'
12	L	89	D2T	CG-CB-SB-CB1
22	a	746	PSU	O4'-C1'-C5-C6
53	V	55	PSU	O4'-C1'-C5-C6
25	d	150	MEQ	CA-CB-CG-CD
55	X	15	7SN	C4'-C5'-O5'-P
22	a	2069	G7M	O4'-C4'-C5'-O5'
22	a	2503	2MA	O4'-C4'-C5'-O5'
53	V	35	U8U	C2'-C1'-N1-C2
22	a	2445	2MG	O4'-C4'-C5'-O5'
33	l	82	MS6	CB-CG-SD-CE
33	l	81	4D4	O-C-CA-CB
55	X	56	OMC	C2'-C1'-N1-C2

There are no ring outliers.

16 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	a	2552	OMU	1	0
22	a	747	5MU	1	0
53	V	55	PSU	2	0
55	X	26	M2G	1	0
22	a	2604	PSU	1	0
53	V	54	5MU	1	0
1	A	1519	MA6	1	0
55	X	15	7SN	1	0
53	V	35	U8U	1	0
11	K	119	IAS	1	0
22	a	1939	5MU	1	0
22	a	2030	6MZ	2	0
1	A	1516	2MG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	a	2251	OMG	1	0
22	a	1917	PSU	1	0
22	a	2503	2MA	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 318 ligands modelled in this entry, 318 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.

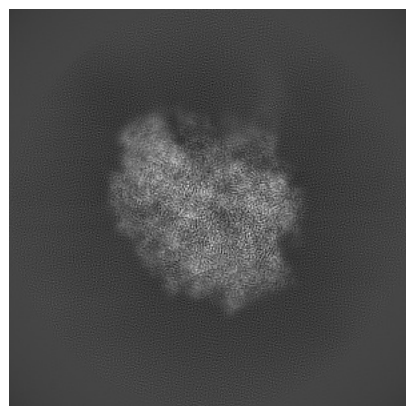
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-35020. These allow visual inspection of the internal detail of the map and identification of artifacts.

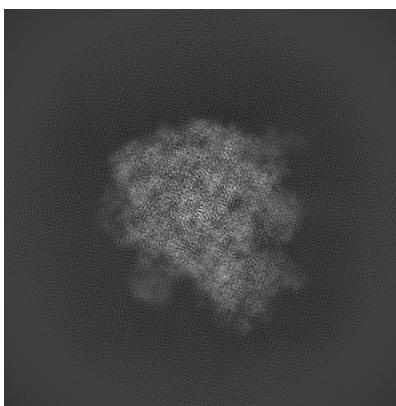
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

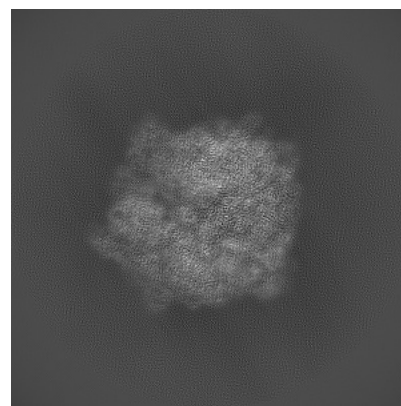
6.1.1 Primary map



X

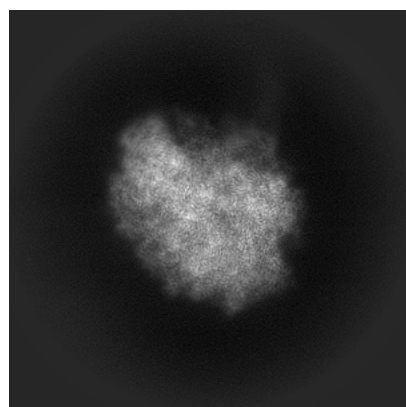


Y

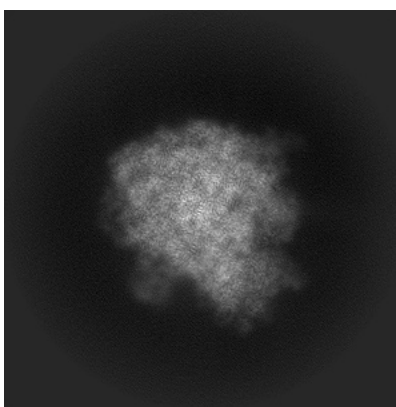


Z

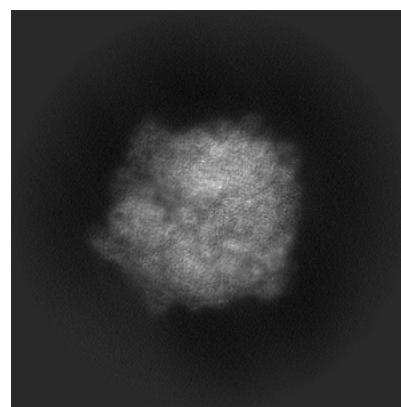
6.1.2 Raw map



X



Y

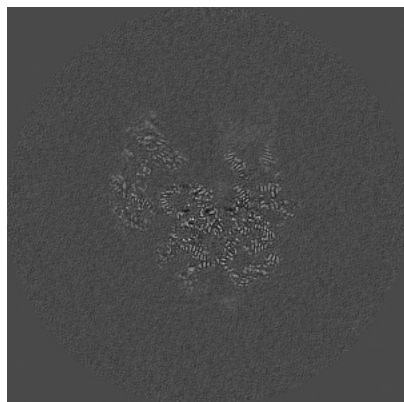


Z

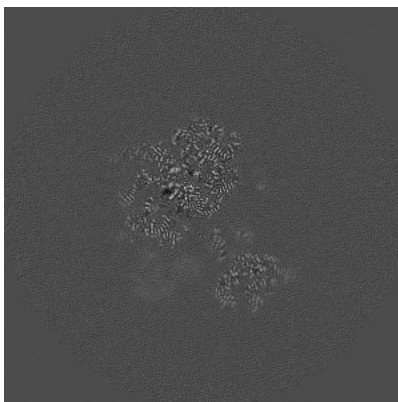
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

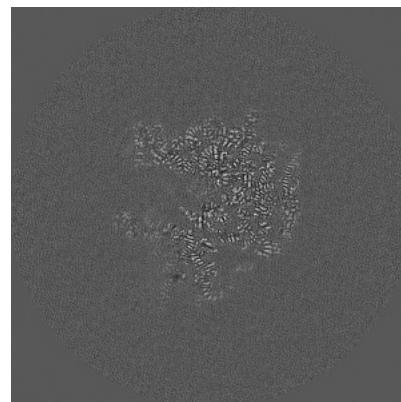
6.2.1 Primary map



X Index: 265

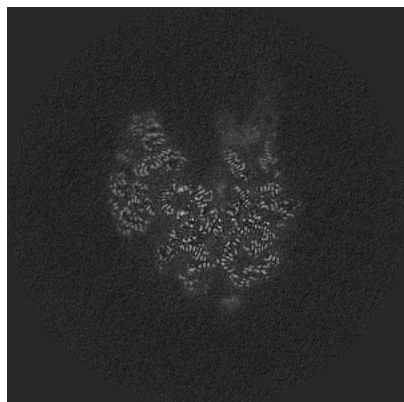


Y Index: 265

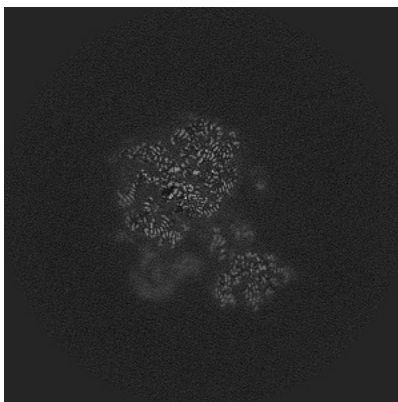


Z Index: 265

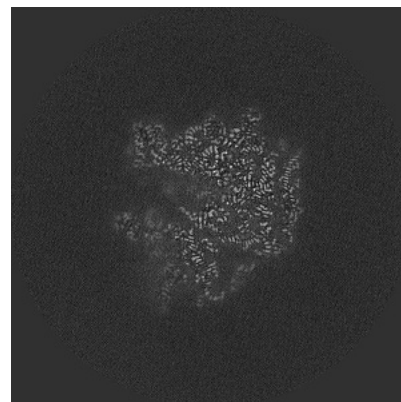
6.2.2 Raw map



X Index: 265



Y Index: 265

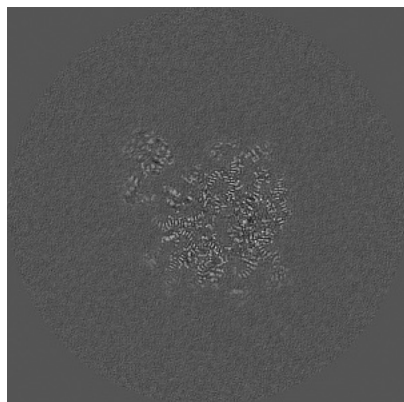


Z Index: 265

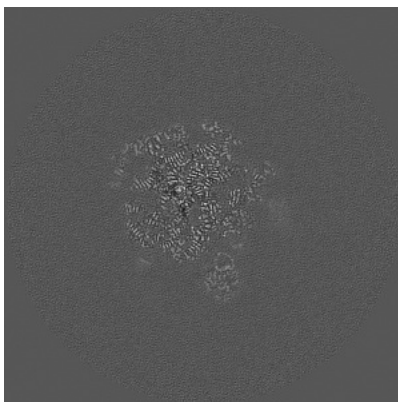
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

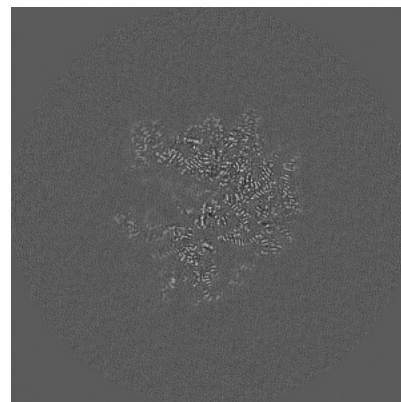
6.3.1 Primary map



X Index: 297

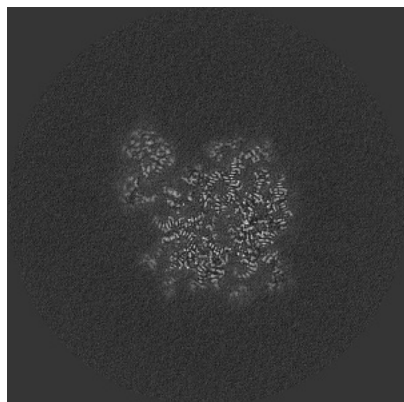


Y Index: 318

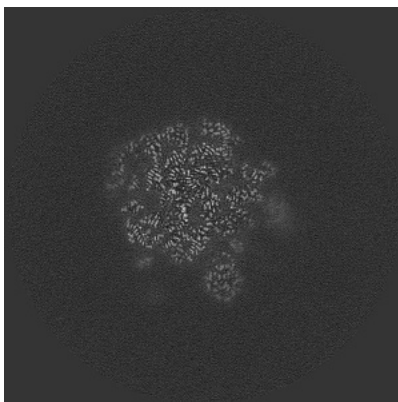


Z Index: 268

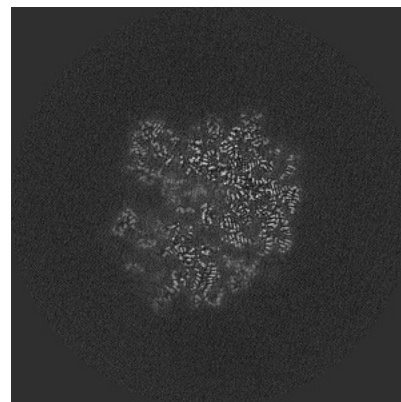
6.3.2 Raw map



X Index: 298



Y Index: 316

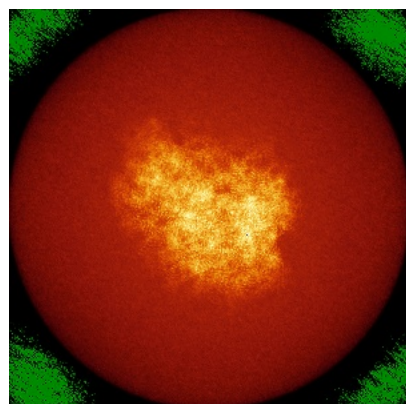


Z Index: 272

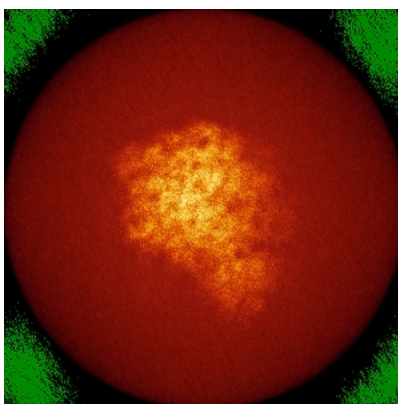
The images above show the largest variance slices of the map in three orthogonal directions.

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

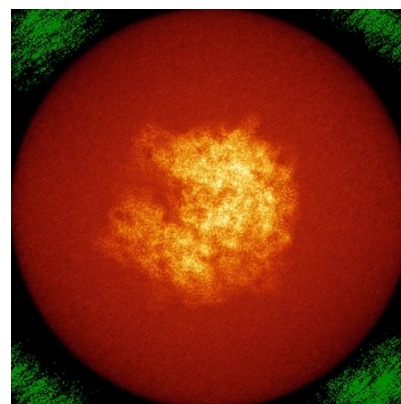
6.4.1 Primary map



X

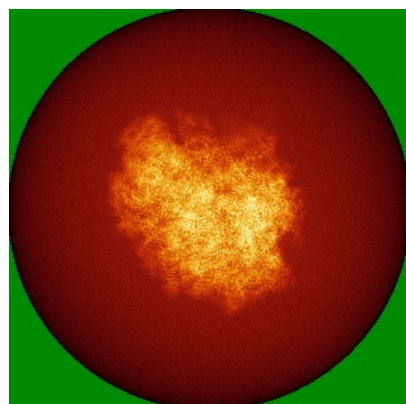


Y

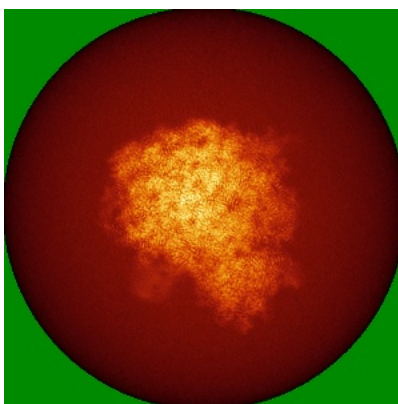


Z

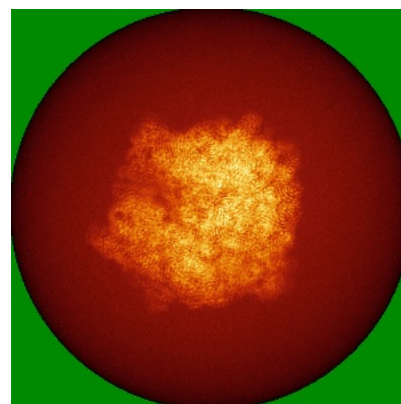
6.4.2 Raw map



X



Y

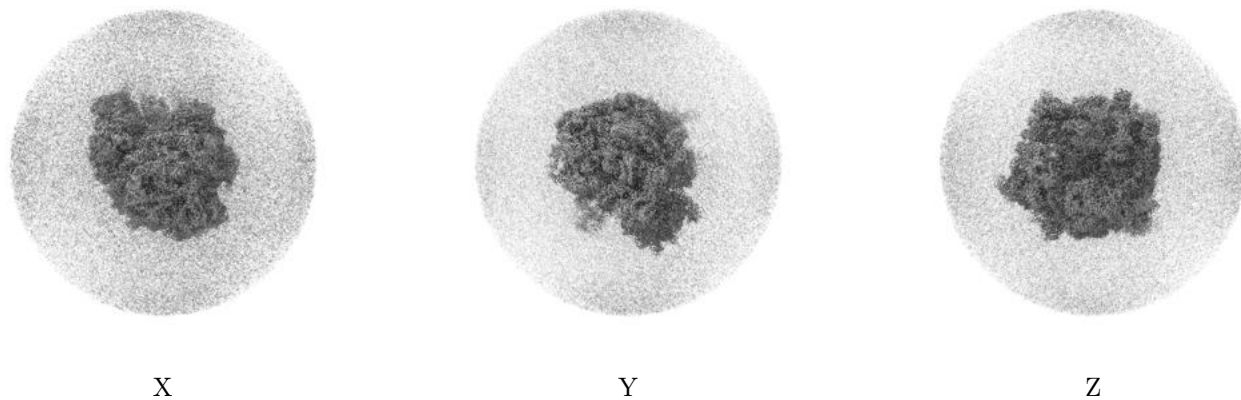


Z

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.

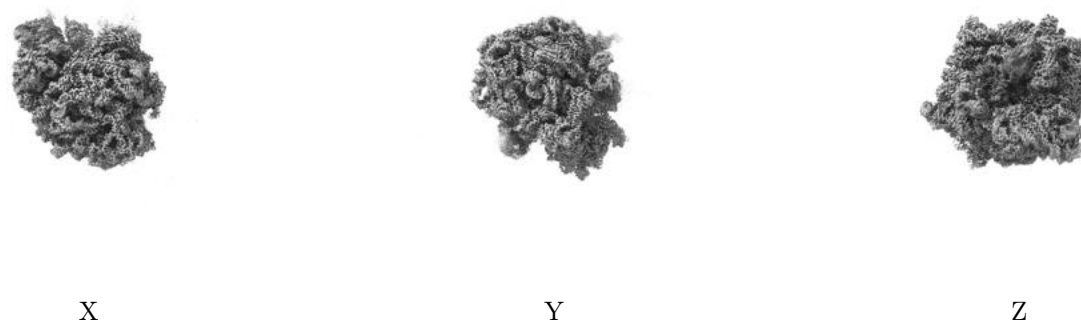
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0253. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

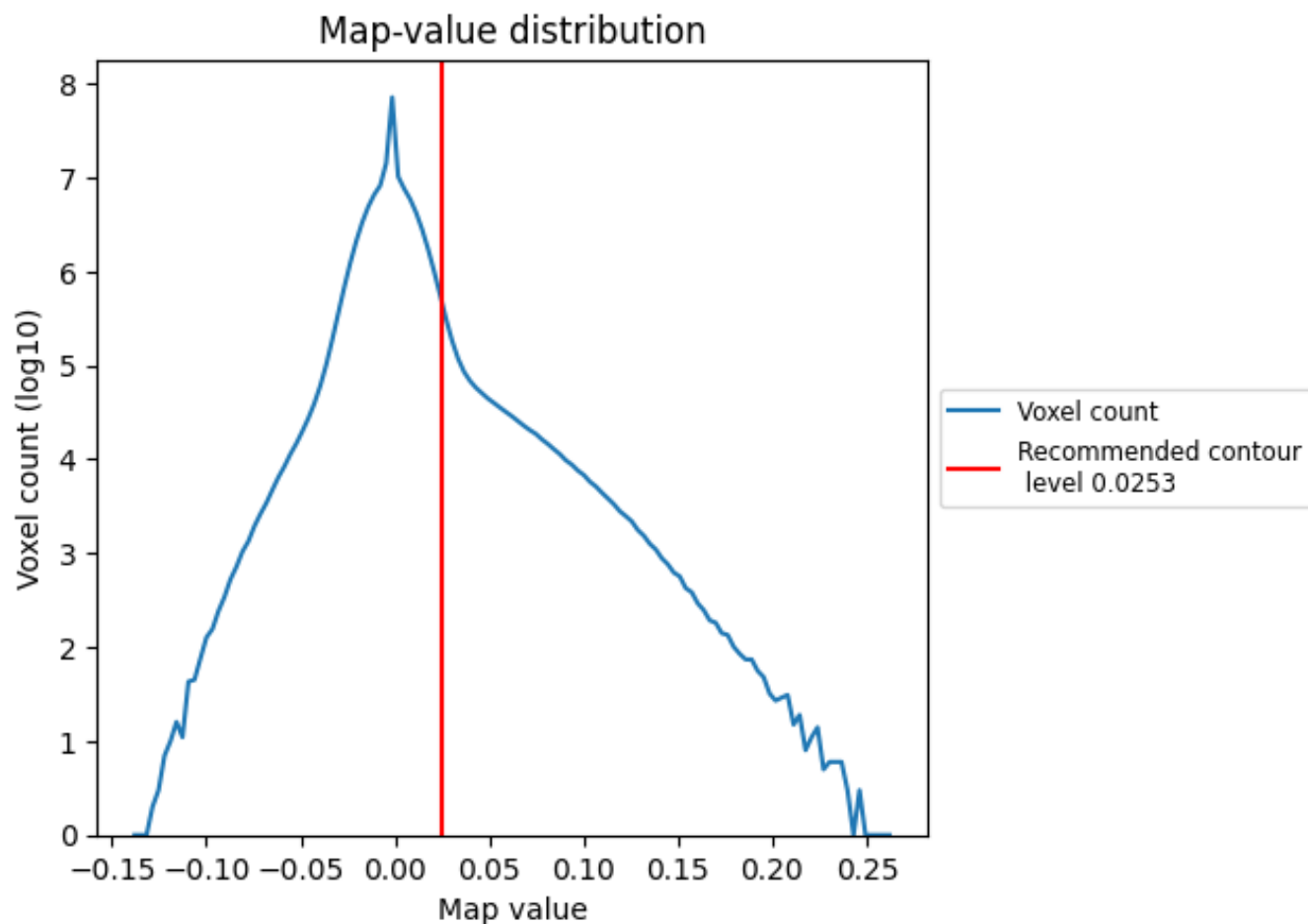
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

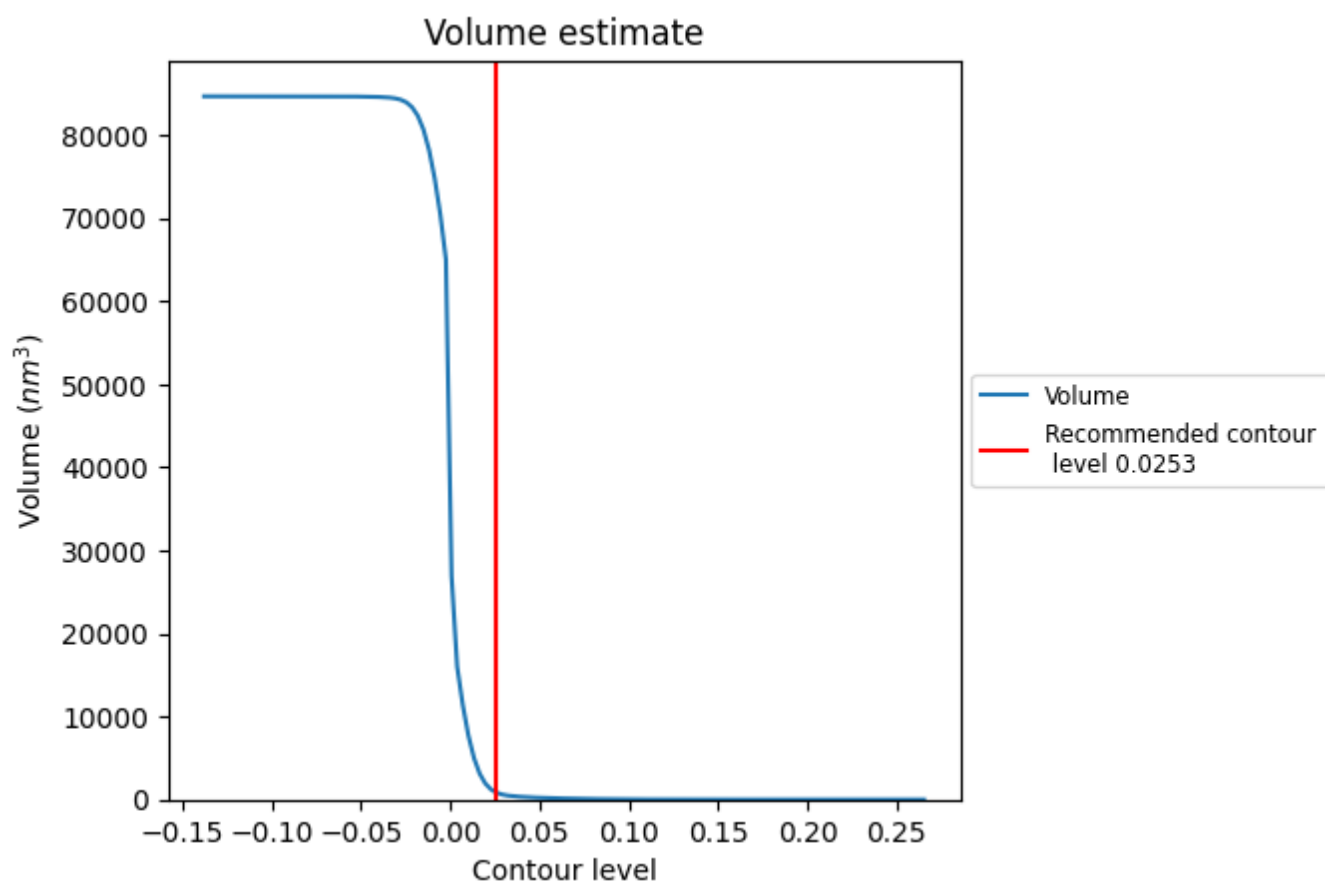
This section contains the results of statistical analysis of the map.

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.

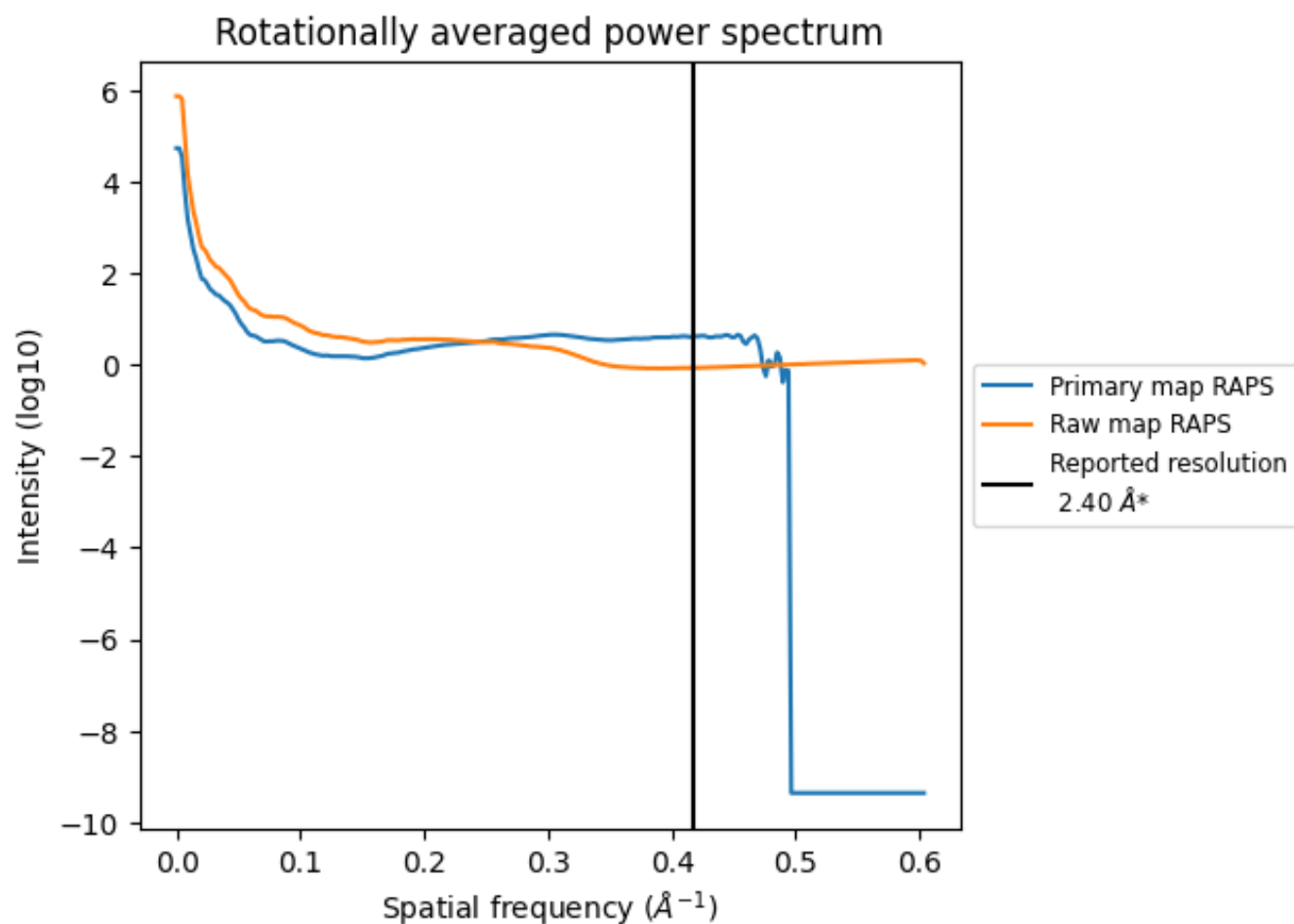
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 902 nm^3 ; this corresponds to an approximate mass of 815 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 ⓘ

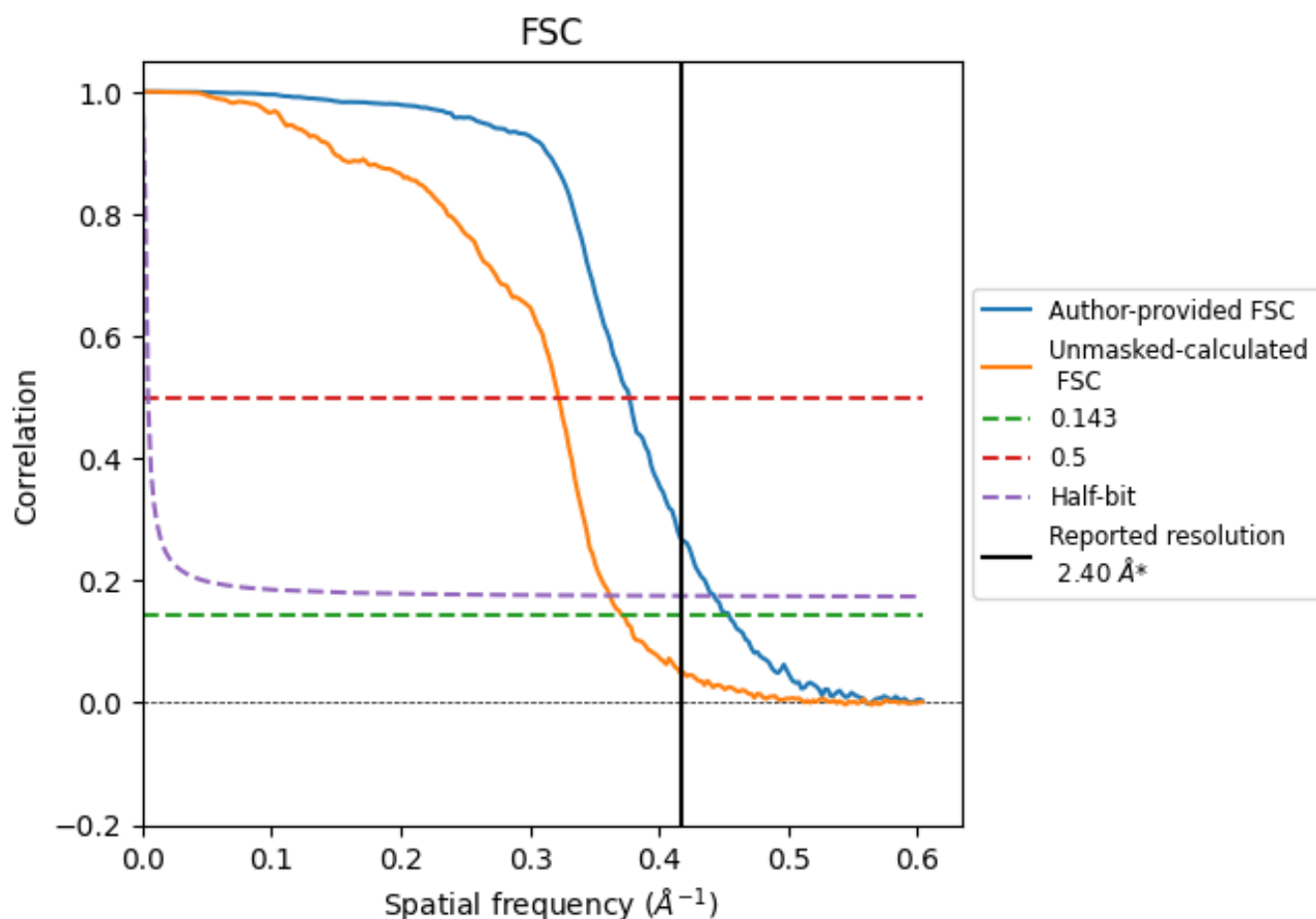


*Reported resolution corresponds to spatial frequency of 0.417 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.417 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

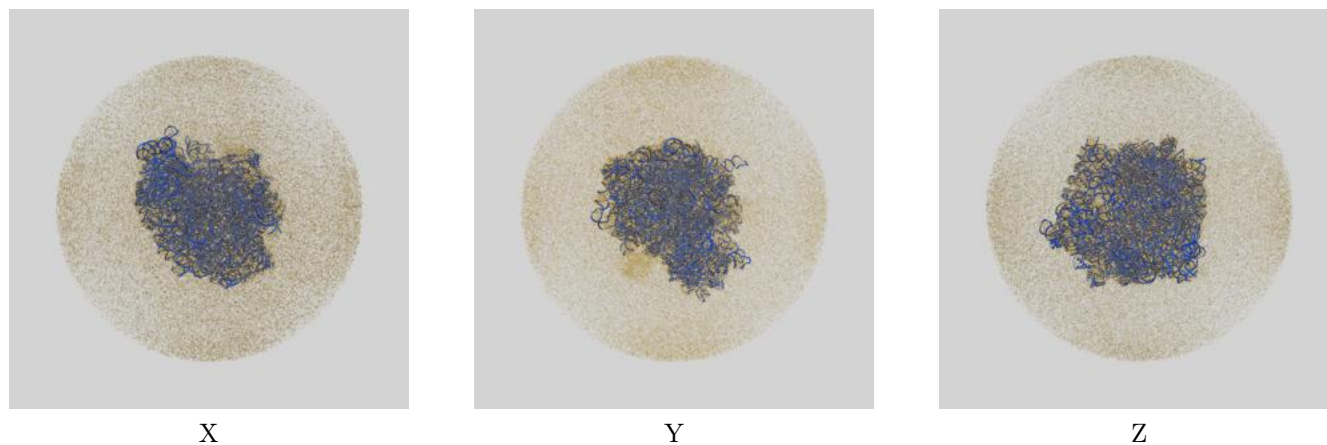
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.40	-	-
Author-provided FSC curve	2.20	2.65	2.26
Unmasked-calculated*	2.69	3.10	2.76

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 2.69 differs from the reported value 2.4 by more than 10 %

9 Map-model fit [i](#)

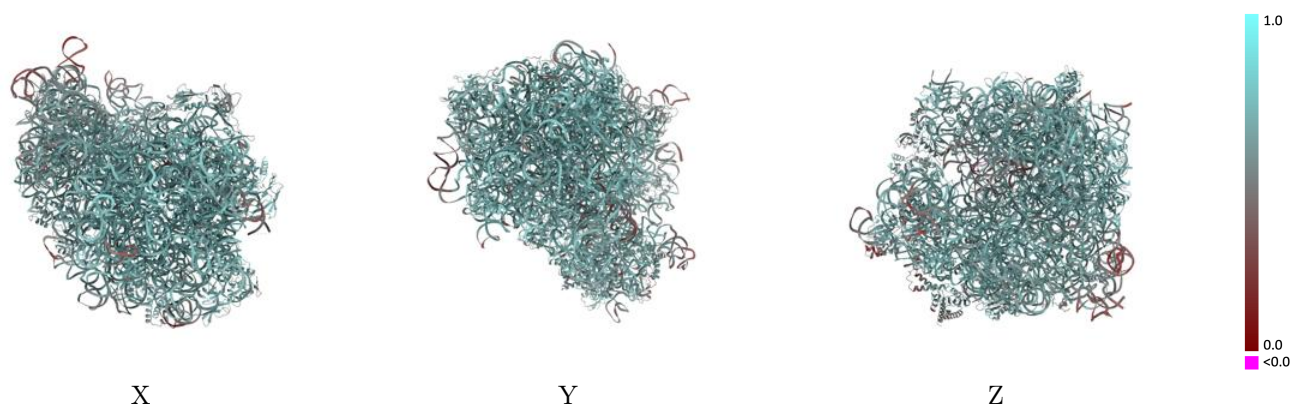
This section contains information regarding the fit between EMDB map EMD-35020 and PDB model 8HTZ. Per-residue inclusion information can be found in section 3 on page 15.

9.1 Map-model overlay [i](#)



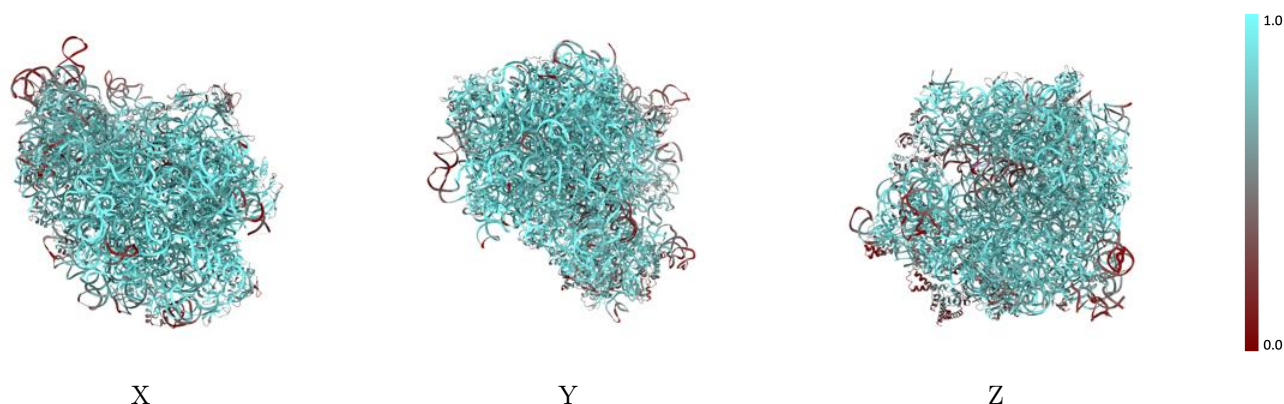
The images above show the 3D surface view of the map at the recommended contour level 0.0253 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



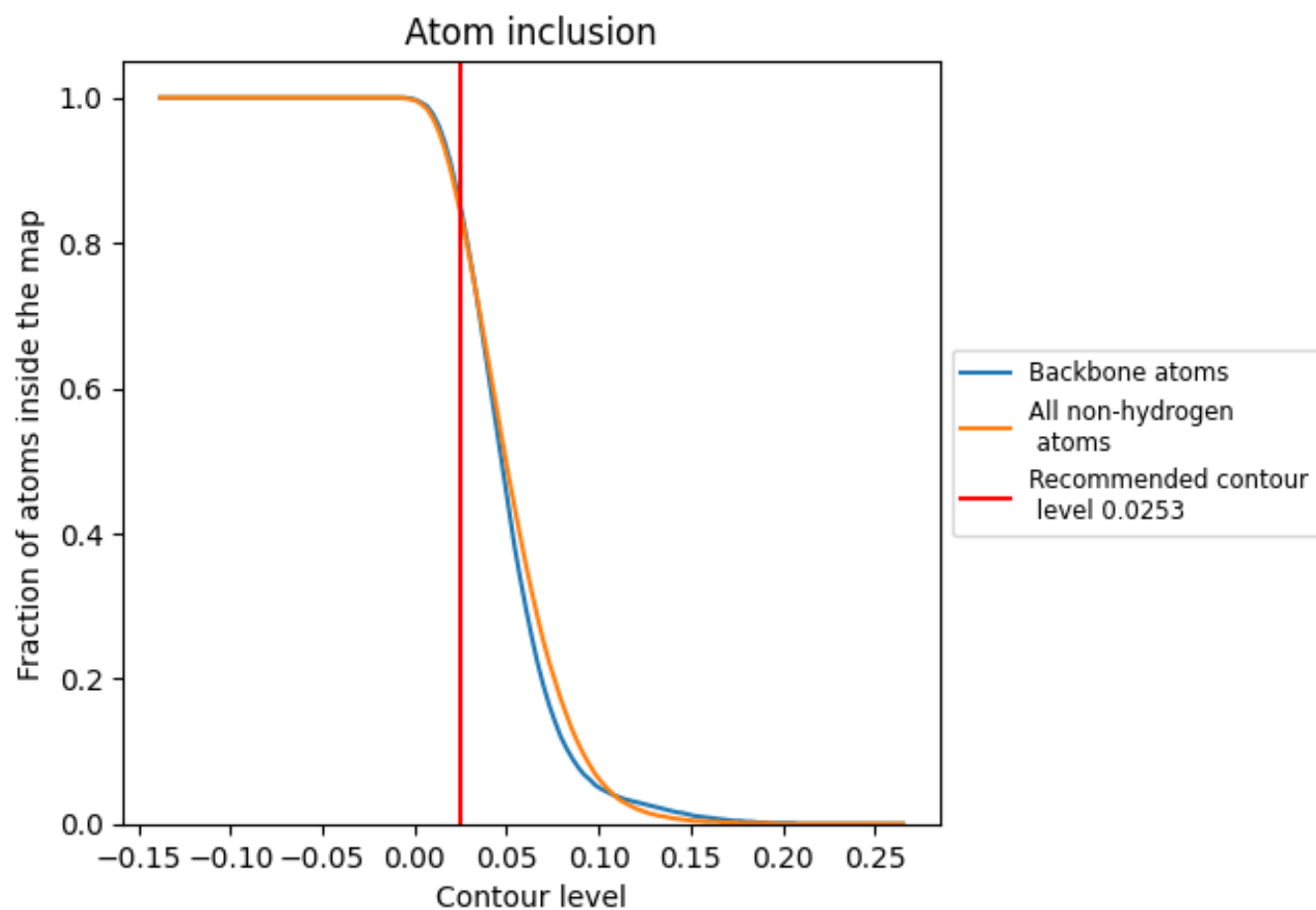
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0253).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 84% 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 (0.0253) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8380	 0.6570
0	 0.8440	 0.6900
1	 0.9580	 0.7290
2	 0.9630	 0.7230
3	 0.8770	 0.6870
4	 0.4370	 0.5300
A	 0.8360	 0.6420
B	 0.4310	 0.5470
C	 0.7760	 0.6500
D	 0.5650	 0.5900
E	 0.8190	 0.6700
F	 0.6370	 0.5990
G	 0.6840	 0.6200
H	 0.8180	 0.6670
I	 0.7630	 0.6390
J	 0.5640	 0.5620
K	 0.7840	 0.6550
L	 0.8420	 0.6820
M	 0.7390	 0.6350
N	 0.8130	 0.6560
O	 0.7880	 0.6430
P	 0.6420	 0.5840
Q	 0.7010	 0.6160
R	 0.7530	 0.6370
S	 0.7360	 0.6280
T	 0.6400	 0.5980
U	 0.4720	 0.5530
V	 0.5990	 0.5470
W	 0.8490	 0.6540
X	 0.4430	 0.4800
a	 0.9090	 0.6810
b	 0.8560	 0.6430
c	 0.9320	 0.7220
d	 0.9080	 0.7120
e	 0.7720	 0.6550



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Chain	Atom inclusion	Q-score
f	 0.6340	 0.5990
g	 0.5680	 0.5700
h	 0.5370	 0.5800
i	 0.9130	 0.7050
j	 0.8710	 0.6940
k	 0.8760	 0.6880
l	 0.9060	 0.7050
m	 0.9470	 0.7190
n	 0.7910	 0.6470
o	 0.8630	 0.6880
p	 0.9520	 0.7220
q	 0.8290	 0.6700
r	 0.8890	 0.6850
s	 0.7770	 0.6460
t	 0.6940	 0.6080
u	 0.7790	 0.6610
v	 0.9060	 0.7100
w	 0.8550	 0.6850
x	 0.6420	 0.5870
y	 0.8630	 0.6760
z	 0.8700	 0.6930