



Full wwPDB EM Validation Report ⓘ

Jun 9, 2026 – 05:02 pm BST

PDB ID : 9SS5 / pdb_00009ss5
EMDB ID : EMD-55177
Title : 50S ribosomal subunit in rimM-KO with RsfS (+H68)
Authors : Hassan, A.H.; Demo, G.
Deposited on : 2025-09-25
Resolution : 2.70 Å (reported)
Based on initial model : 6WDE

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:

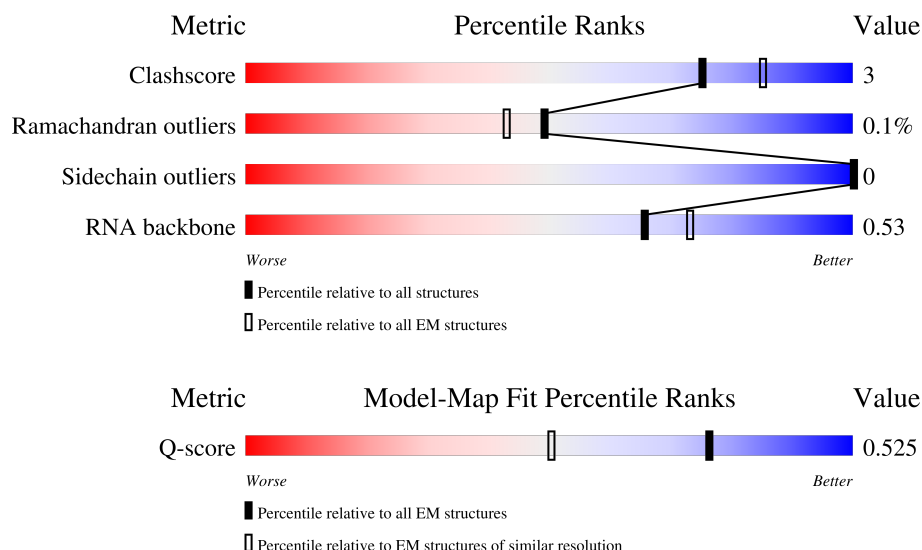
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 2.70 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.







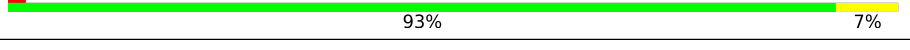

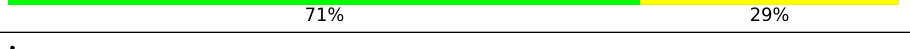
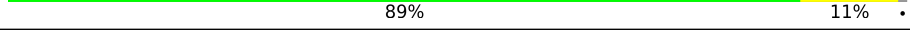
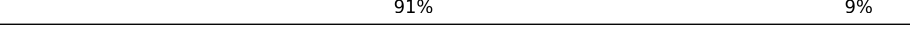
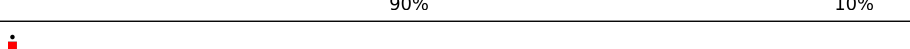
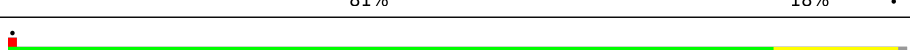

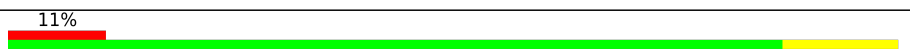

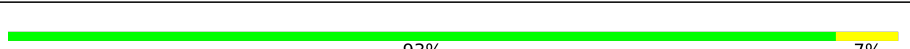

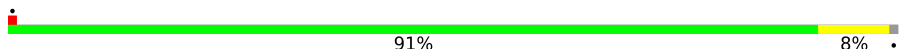



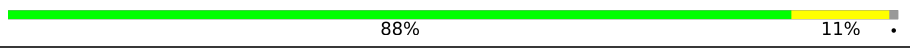
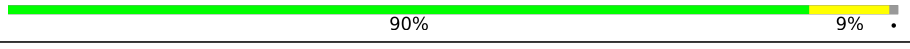



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	10327 (2.20 - 3.20)

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	B0	85	
2	B1	78	
3	B2	63	

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Mol	Chain	Length	Quality of chain
4	B3	59	
5	B4	70	
6	B5	57	
7	B6	55	
8	B7	46	
9	B8	65	
10	B9	38	
11	BB	273	
12	BC	209	
13	BD	201	
14	BE	179	
15	BF	177	
16	BG	87	
17	BI	149	
18	BK	142	
19	BM	142	
20	BN	123	
21	BO	144	
22	BP	136	
23	BQ	127	
24	BR	117	
25	BS	115	
26	BT	118	
27	BU	103	
28	BV	110	

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Mol	Chain	Length	Quality of chain
29	BW	100	
30	BX	104	
31	BY	94	
32	D2	2903	
33	D3	120	
34	FS	105	

2 Entry composition

There are 34 unique types of molecules in this entry. The entry contains 92714 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 protein called Large ribosomal subunit protein bL27.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B0	75	Total	C	N	O	S	0	0
			575	356	116	102	1		

- Molecule 2 is a protein called Large ribosomal subunit protein bL28.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B1	77	Total	C	N	O	S	0	0
			625	388	129	106	2		

- Molecule 3 is a protein called Large ribosomal subunit protein uL29.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B2	63	Total	C	N	O	S	0	0
			509	313	99	95	2		

- Molecule 4 is a protein called Large ribosomal subunit protein uL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	B3	58	Total	C	N	O	S	0	0
			449	281	87	79	2		

- Molecule 5 is a protein called Large ribosomal subunit protein bL31A.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	B4	47	Total	C	N	O	S	0	0
			364	227	64	67	6		

- Molecule 6 is a protein called Large ribosomal subunit protein bL32.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	B5	56	Total	C	N	O	S	0	0
			444	269	94	80	1		

- Molecule 7 is a protein called Large ribosomal subunit protein bL33.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	B6	50	Total	C	N	O	0	0
			409	263	75	71		

- Molecule 8 is a protein called Large ribosomal subunit protein bL34.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	B7	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 9 is a protein called Large ribosomal subunit protein bL35.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	B8	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 10 is a protein called Large ribosomal subunit protein bL36A.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	B9	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 11 is a protein called Large ribosomal subunit protein uL2.

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

- Molecule 12 is a protein called 50S ribosomal protein L3.

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

- Molecule 13 is a protein called Large ribosomal subunit protein uL4.

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

- Molecule 14 is a protein called Large ribosomal subunit protein uL5.

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

- Molecule 15 is a protein called Large ribosomal subunit protein uL6.

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

- Molecule 16 is a protein called Small ribosomal subunit protein bS20.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	BG	71	Total	C	N	O	S	0	0
			557	346	115	93	3		

- Molecule 17 is a protein called Large ribosomal subunit protein bL9.

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

- Molecule 18 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	BK	141	Total	C	N	O	S	0	0
			1032	651	179	196	6		

- Molecule 19 is a protein called Large ribosomal subunit protein uL13.

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

- Molecule 20 is a protein called Large ribosomal subunit protein uL14.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	BN	122	Total	C	N	O	S	0	0
			938	587	180	165	6		

- Molecule 21 is a protein called Large ribosomal subunit protein uL15.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	BO	143	Total	C	N	O	S	0	0
			1045	649	206	189	1		

- Molecule 22 is a protein called 50S ribosomal protein L16.

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

- Molecule 23 is a protein called Large ribosomal subunit protein bL17.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	BQ	120	Total	C	N	O	S	0	0
			960	593	196	166	5		

- Molecule 24 is a protein called Large ribosomal subunit protein uL18.

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

- Molecule 25 is a protein called Large ribosomal subunit protein bL19.

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

- Molecule 26 is a protein called Large ribosomal subunit protein bL20.

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

- Molecule 27 is a protein called Large ribosomal subunit protein bL21.

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

- Molecule 28 is a protein called Large ribosomal subunit protein uL22.

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

- Molecule 29 is a protein called Large ribosomal subunit protein uL23.

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

- Molecule 30 is a protein called Large ribosomal subunit protein uL24.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	BX	102	Total	C	N	O		0	0
			779	492	146	141			

- Molecule 31 is a protein called Large ribosomal subunit protein bL25.

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

- Molecule 32 is a RNA chain called 23S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	D2	2903	Total	C	N	O	P	0	0
			62317	27801	11468	20146	2902		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D2	747	C	U	conflict	GB CP194060.1

- Molecule 33 is a RNA chain called 5S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	D3	120	Total	C	N	O	P	0	0
			2568	1145	471	833	119		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D3	120	A	U	conflict	GB CP076697.1


- Molecule 34 is a protein called Ribosomal silencing factor RsfS.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
34	FS	103	794	496	135	158	5	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and 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: Large ribosomal subunit protein bL27

Chain B0: 



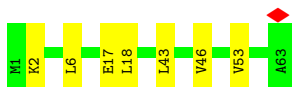
- Molecule 2: Large ribosomal subunit protein bL28

Chain B1: 




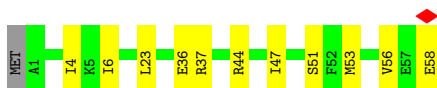
- Molecule 3: Large ribosomal subunit protein uL29

Chain B2: 



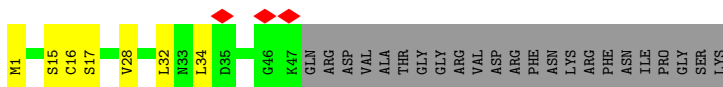
- Molecule 4: Large ribosomal subunit protein uL30

Chain B3: 



- Molecule 5: Large ribosomal subunit protein bL31A

Chain B4: 




- Molecule 6: Large ribosomal subunit protein bL32

Chain B5:  89% 9%



- Molecule 7: Large ribosomal subunit protein bL33

Chain B6:  78% 13% 9%




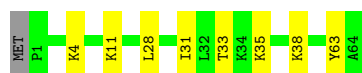
- Molecule 8: Large ribosomal subunit protein bL34

Chain B7:  93% 7%



- Molecule 9: Large ribosomal subunit protein bL35

Chain B8:  86% 12%




- Molecule 10: Large ribosomal subunit protein bL36A

Chain B9:  71% 29%



- Molecule 11: Large ribosomal subunit protein uL2

Chain BB:  89% 11%



- Molecule 12: 50S ribosomal protein L3

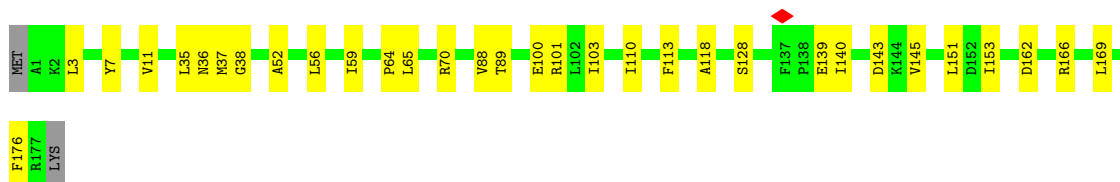
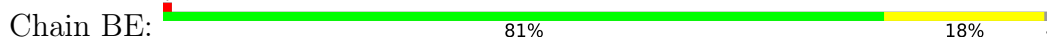
Chain BC:  91% 9%



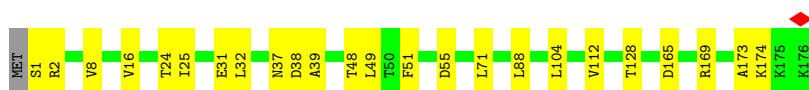
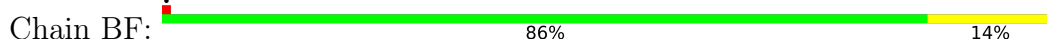
- Molecule 13: Large ribosomal subunit protein uL4



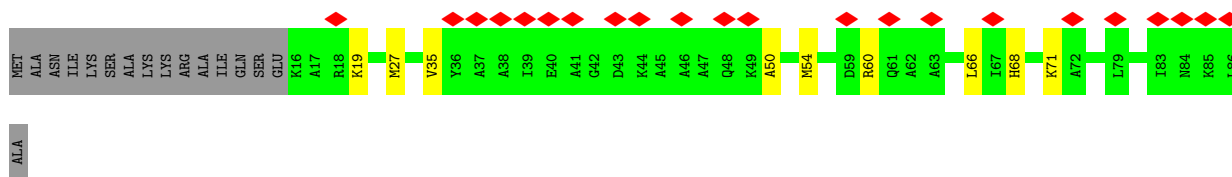
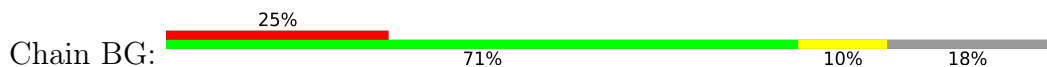
- Molecule 14: Large ribosomal subunit protein uL5



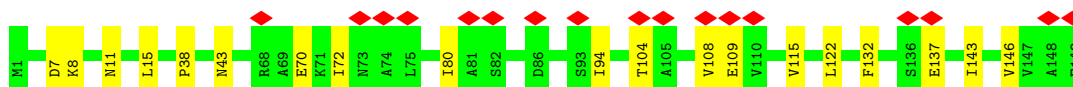
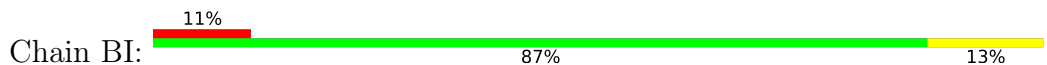
- Molecule 15: Large ribosomal subunit protein uL6



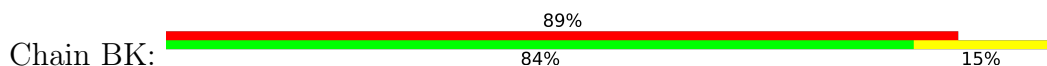
- Molecule 16: Small ribosomal subunit protein bS20

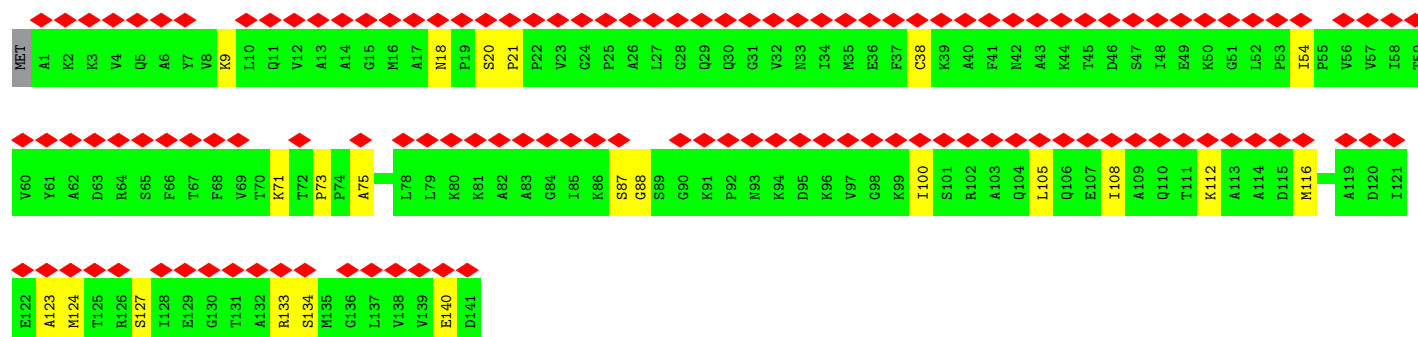


- Molecule 17: Large ribosomal subunit protein bL9



- Molecule 18: 50S ribosomal protein L11





• Molecule 19: Large ribosomal subunit protein uL13

Chain BM: 93% 7%



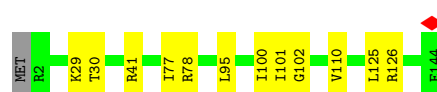
• Molecule 20: Large ribosomal subunit protein uL14

Chain BN: 83% 16%



• Molecule 21: Large ribosomal subunit protein uL15

Chain BO: 91% 8%



• Molecule 22: 50S ribosomal protein L16

Chain BP: 81% 19%



• Molecule 23: Large ribosomal subunit protein bL17

Chain BQ: 83% 11% 6%



• Molecule 24: Large ribosomal subunit protein uL18

Chain BR: 85% 15% 6%



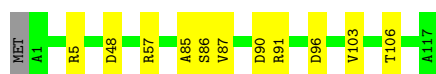
- Molecule 25: Large ribosomal subunit protein bL19

Chain BS: 88% 11% .



- Molecule 26: Large ribosomal subunit protein bL20

Chain BT: 90% 9% .



- Molecule 27: Large ribosomal subunit protein bL21

Chain BU: 85% 14% .



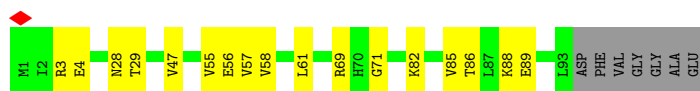
- Molecule 28: Large ribosomal subunit protein uL22

Chain BV: 85% 15%



- Molecule 29: Large ribosomal subunit protein uL23

Chain BW: 76% 17% 7%



- Molecule 30: Large ribosomal subunit protein uL24

Chain BX: 89% 8% ..



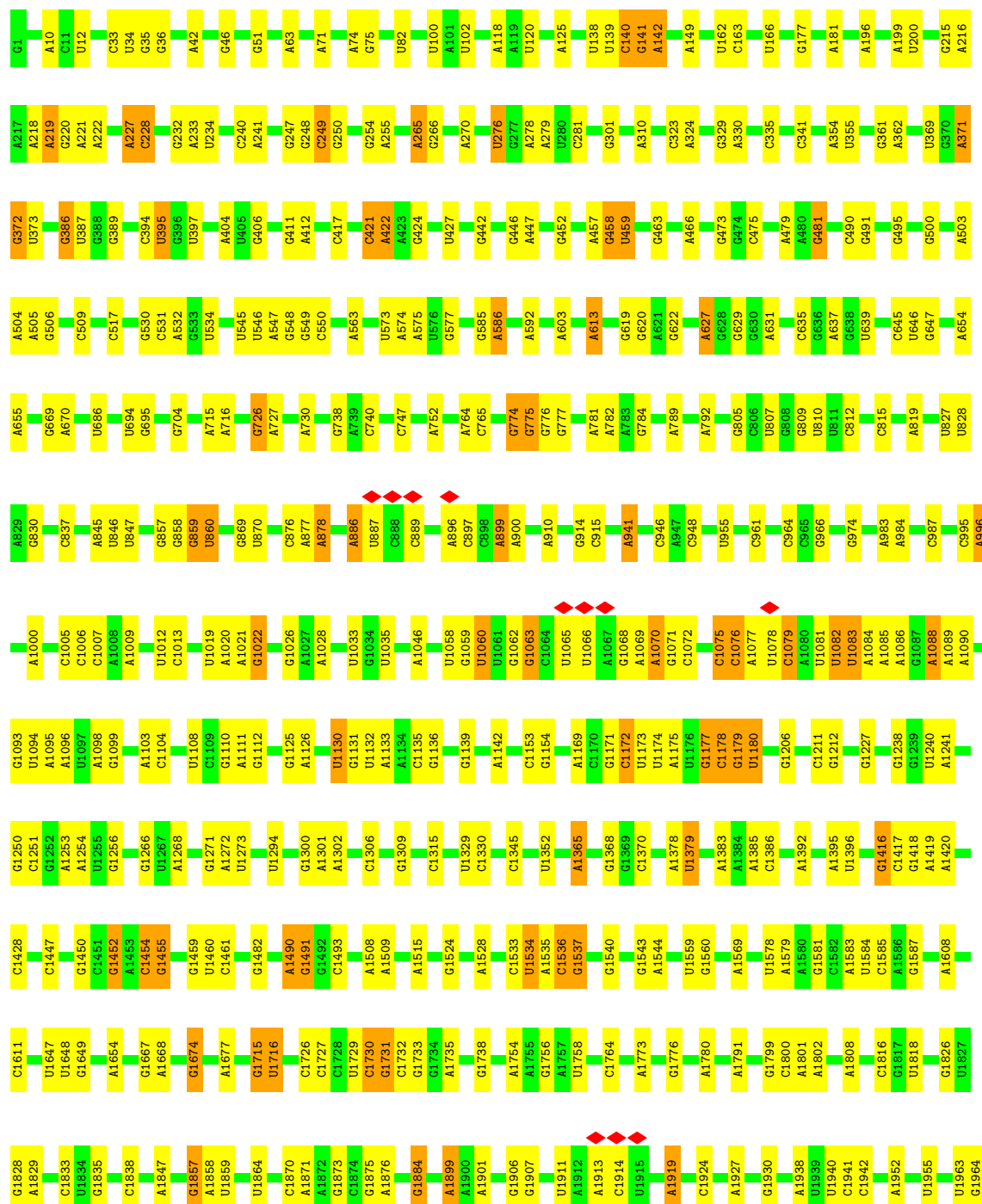
- Molecule 31: Large ribosomal subunit protein bL25

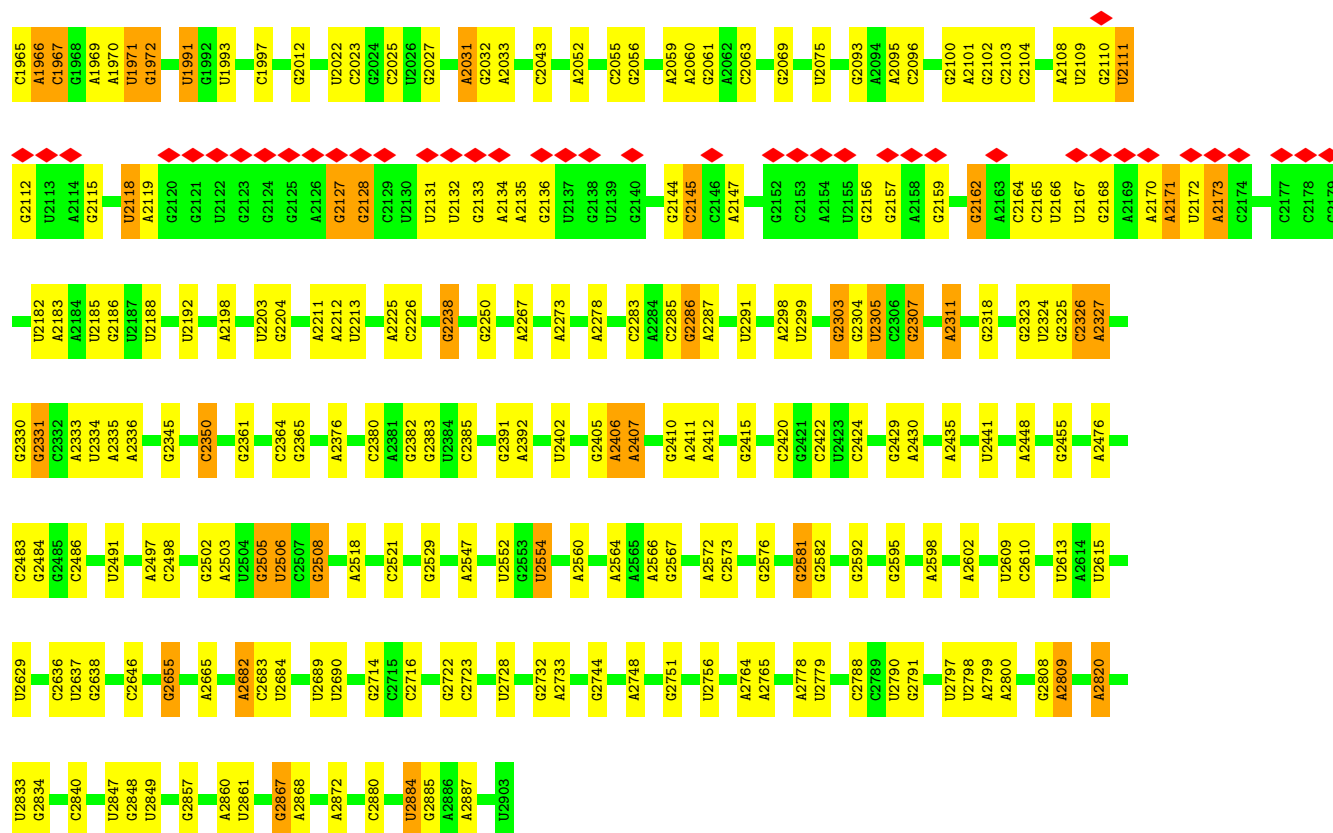
Chain BY: 96% .



• Molecule 32: 23S Ribosomal RNA

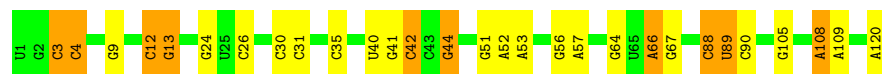
Chain D2: 77% 20%





• Molecule 33: 5S Ribosomal RNA

Chain D3: 76% 16% 8%



• Molecule 34: Ribosomal silencing factor RsfS

Chain FS: 82% 16% 2%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	51537	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	22.078	Depositor
Minimum map value	-6.685	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.972	Depositor
Recommended contour level	2.2	Depositor
Map size (Å)	400.32, 400.32, 400.32	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.834, 0.834, 0.834	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	B0	0.18	0/582	0.31	0/769
2	B1	0.19	0/635	0.29	0/848
3	B2	0.17	0/510	0.42	0/677
4	B3	0.19	0/453	0.35	0/605
5	B4	0.15	0/371	0.36	0/496
6	B5	0.20	0/450	0.38	0/599
7	B6	0.16	0/416	0.35	0/554
8	B7	0.21	0/380	0.32	0/498
9	B8	0.17	0/513	0.33	0/676
10	B9	0.23	0/303	0.47	0/397
11	BB	0.18	0/2121	0.35	0/2852
12	BC	0.19	0/1586	0.33	0/2134
13	BD	0.17	0/1571	0.33	0/2113
14	BE	0.17	0/1434	0.39	0/1926
15	BF	0.17	0/1343	0.39	0/1816
16	BG	0.16	0/563	0.35	0/747
17	BI	0.16	0/1122	0.38	0/1515
18	BK	0.15	0/1046	0.34	0/1410
19	BM	0.18	0/1152	0.28	0/1551
20	BN	0.18	0/947	0.37	0/1268
21	BO	0.18	0/1054	0.39	0/1403
22	BP	0.19	0/1093	0.42	0/1460
23	BQ	0.20	0/973	0.42	0/1301
24	BR	0.16	0/902	0.38	0/1209
25	BS	0.20	0/929	0.38	0/1242
26	BT	0.22	0/960	0.34	0/1278
27	BU	0.20	0/829	0.45	0/1107
28	BV	0.19	0/864	0.37	0/1156
29	BW	0.19	0/744	0.35	0/994
30	BX	0.19	0/787	0.46	2/1051 (0.2%)
31	BY	0.18	0/766	0.36	0/1025
32	D2	0.20	0/69796	0.29	0/108888
33	D3	0.16	0/2872	0.27	0/4479
34	FS	0.17	0/803	0.41	0/1085

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.19	0/100870	0.31	2/151129 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
30	BX	97	SER	CA-C-N	6.02	133.03	121.54
30	BX	97	SER	C-N-CA	6.02	133.03	121.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B0	575	0	592	9	0
2	B1	625	0	655	7	0
3	B2	509	0	543	4	0
4	B3	449	0	491	8	0
5	B4	364	0	366	6	0
6	B5	444	0	461	3	0
7	B6	409	0	440	7	0
8	B7	377	0	418	2	0
9	B8	504	0	574	7	0
10	B9	302	0	343	8	0
11	BB	2082	0	2157	19	0
12	BC	1565	0	1616	12	0
13	BD	1552	0	1619	14	0
14	BE	1410	0	1447	25	0
15	BF	1323	0	1374	18	0
16	BG	557	0	595	7	0
17	BI	1111	0	1148	13	0
18	BK	1032	0	1088	18	0
19	BM	1129	0	1162	9	0
20	BN	938	0	1012	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
21	BO	1045	0	1117	9	0
22	BP	1074	0	1157	16	0
23	BQ	960	0	1000	10	0
24	BR	892	0	923	10	0
25	BS	917	0	965	8	0
26	BT	947	0	1022	9	0
27	BU	816	0	839	11	0
28	BV	857	0	922	10	0
29	BW	738	0	807	11	0
30	BX	779	0	834	6	0
31	BY	753	0	780	2	0
32	D2	62317	0	31346	221	0
33	D3	2568	0	1303	13	0
34	FS	794	0	793	12	0
All	All	92714	0	61909	465	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 3.

All (465) 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)
32:D2:1173:U:O2'	32:D2:1177:G:N2	2.05	0.90
32:D2:2848:G:O2'	32:D2:2867:G:N2	2.09	0.86
14:BE:35:LEU:HD12	14:BE:151:LEU:HD21	1.59	0.84
32:D2:2326:C:O2'	32:D2:2327:A:OP1	1.93	0.84
22:BP:58:LYS:O	22:BP:59:ARG:NE	2.11	0.83
26:BT:48:ASP:OD2	32:D2:534:U:O2'	1.95	0.83
26:BT:91:ARG:NH2	32:D2:1153:C:OP1	2.13	0.81
32:D2:2406:A:OP2	32:D2:2411:A:N6	2.15	0.80
32:D2:475:C:O2	32:D2:479:A:N6	2.13	0.80
32:D2:270:A:N1	32:D2:369:U:O2'	2.15	0.79
16:BG:68:HIS:NE2	32:D2:42:A:OP1	2.16	0.79
23:BQ:56:LYS:NZ	23:BQ:87:PHE:O	2.14	0.78
12:BC:33:ARG:NH1	12:BC:74:GLU:OE2	2.16	0.78
5:B4:1:MET:N	33:D3:40:U:O4	2.17	0.78
32:D2:704:G:O2'	32:D2:727:A:N6	2.16	0.78
32:D2:1940:U:O2	32:D2:1942:C:N4	2.17	0.77
23:BQ:69:ARG:O	23:BQ:70:THR:OG1	2.03	0.77
32:D2:200:U:O2	32:D2:386:G:N2	2.18	0.77
22:BP:57:VAL:O	22:BP:60:GLN:NE2	2.18	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:D2:2118:U:O2	32:D2:2145:C:N4	2.18	0.75
24:BR:111:ARG:NH1	24:BR:117:PHE:OXT	2.20	0.75
14:BE:59:ILE:O	14:BE:101:ARG:NH2	2.19	0.75
17:BI:38:PRO:O	17:BI:43:ASN:ND2	2.20	0.75
20:BN:9:ASN:OD1	20:BN:18:ARG:NH1	2.19	0.75
28:BV:22:ASP:OD1	28:BV:25:ARG:NH1	2.21	0.74
29:BW:28:ASN:ND2	29:BW:88:LYS:O	2.20	0.74
26:BT:57:ARG:NH1	32:D2:1154:G:OP2	2.21	0.73
32:D2:220:G:O2'	32:D2:233:A:N3	2.21	0.73
32:D2:1911:U:O2	32:D2:1919:A:N6	2.21	0.73
32:D2:1858:A:N6	32:D2:1884:G:O2'	2.20	0.73
21:BO:29:LYS:O	21:BO:30:THR:OG1	2.05	0.73
20:BN:70:ARG:NH2	32:D2:2683:C:O2	2.21	0.73
1:B0:74:LYS:NZ	32:D2:858:G:OP1	2.21	0.73
10:B9:30:GLU:N	10:B9:30:GLU:OE1	2.22	0.73
4:B3:6:ILE:HD13	4:B3:47:ILE:HD11	1.69	0.72
32:D2:2646:C:OP2	32:D2:2732:G:O2'	2.08	0.72
2:B1:71:ARG:NH1	2:B1:77:TYR:OH	2.22	0.72
23:BQ:106:ASP:OD2	32:D2:1649:G:O2'	2.03	0.72
32:D2:1169:A:N6	32:D2:1180:U:O4	2.14	0.72
32:D2:2127:G:O2'	32:D2:2128:G:O5'	2.07	0.72
32:D2:2581:G:O2'	32:D2:2610:C:N4	2.22	0.72
30:BX:81:ARG:NH2	32:D2:301:G:OP2	2.22	0.71
13:BD:144:GLU:N	13:BD:144:GLU:OE1	2.24	0.71
32:D2:481:G:O2'	32:D2:506:G:N2	2.23	0.71
18:BK:116:MET:SD	32:D2:1058:U:O2'	2.49	0.71
23:BQ:49:GLU:OE2	23:BQ:95:THR:OG1	2.08	0.71
33:D3:88:C:O2'	33:D3:89:U:O5'	2.06	0.70
10:B9:19:ARG:NE	32:D2:2756:U:OP2	2.23	0.70
23:BQ:53:THR:HG21	32:D2:2840:C:H5''	1.73	0.70
22:BP:6:ARG:NH1	32:D2:870:U:OP1	2.24	0.70
32:D2:1490:A:O2'	32:D2:1491:G:OP1	2.10	0.70
32:D2:1864:U:OP1	32:D2:2410:G:O2'	2.10	0.70
19:BM:68:LYS:NZ	32:D2:1022:G:O6	2.19	0.69
32:D2:1093:G:N2	32:D2:1099:G:O6	2.26	0.69
29:BW:3:ARG:NH1	32:D2:138:U:O4	2.26	0.68
32:D2:2307:G:N7	32:D2:2311:A:N6	2.40	0.68
12:BC:77:ARG:NH1	12:BC:200:ASP:OD1	2.25	0.68
21:BO:78:ARG:NH2	32:D2:627:A:OP1	2.27	0.68
32:D2:1447:C:O2'	32:D2:1544:A:N3	2.26	0.68
19:BM:30:THR:HG21	32:D2:1005:C:O2'	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:D2:227:A:O2'	32:D2:228:C:O5'	2.11	0.68
34:FS:14:ILE:HD11	34:FS:40:CYS:HB2	1.76	0.68
7:B6:5:ARG:NH1	32:D2:2285:C:OP2	2.26	0.68
14:BE:128:SER:OG	32:D2:2303:G:O2'	2.11	0.67
28:BV:25:ARG:NH2	28:BV:74:ILE:O	2.27	0.67
10:B9:7:VAL:HG11	10:B9:36:ARG:O	1.94	0.67
14:BE:151:LEU:N	32:D2:2305:U:O4	2.28	0.67
22:BP:86:LYS:NZ	32:D2:955:U:OP1	2.24	0.67
32:D2:276:U:O2'	32:D2:278:A:N6	2.27	0.67
17:BI:11:ASN:ND2	32:D2:2095:A:OP1	2.28	0.67
32:D2:629:G:N3	32:D2:639:U:O2'	2.29	0.66
2:B1:9:LYS:NZ	32:D2:397:U:OP2	2.19	0.66
9:B8:31:ILE:O	9:B8:35:LYS:NZ	2.28	0.66
33:D3:12:C:O2'	33:D3:13:G:O5'	2.13	0.66
13:BD:117:ARG:NH2	13:BD:183:PHE:O	2.28	0.65
4:B3:36:GLU:O	4:B3:37:ARG:NH1	2.30	0.65
6:B5:12:ARG:NH2	32:D2:517:C:OP1	2.30	0.65
11:BB:145:MET:HE3	11:BB:153:LEU:HD11	1.79	0.65
32:D2:2165:C:N4	32:D2:2167:U:O4	2.29	0.65
22:BP:76:LYS:NZ	22:BP:83:GLY:O	2.30	0.65
7:B6:22:THR:HG22	7:B6:23:THR:H	1.60	0.65
20:BN:70:ARG:NH1	32:D2:2684:U:O4'	2.30	0.65
22:BP:74:THR:HG22	22:BP:89:VAL:HA	1.79	0.65
5:B4:16:CYS:SG	5:B4:17:SER:N	2.69	0.65
32:D2:1077:A:N6	32:D2:1089:A:OP1	2.30	0.64
30:BX:9:GLU:OE1	30:BX:21:ARG:NH2	2.31	0.64
10:B9:29:ALA:O	15:BF:169:ARG:NH2	2.30	0.64
11:BB:106:PRO:HD2	11:BB:109:LEU:HD22	1.79	0.64
32:D2:948:C:O2	32:D2:984:A:O2'	2.16	0.64
20:BN:44:LYS:O	20:BN:54:LYS:NZ	2.31	0.64
18:BK:71:LYS:NZ	32:D2:1059:G:OP1	2.26	0.63
7:B6:22:THR:HG21	32:D2:2286:G:O6	1.99	0.63
26:BT:90:ASP:OD2	32:D2:996:A:O2'	2.09	0.63
26:BT:96:ASP:OD2	27:BU:13:ARG:NE	2.32	0.63
32:D2:1266:G:O2'	32:D2:2012:G:O6	2.14	0.62
32:D2:1536:C:HO2'	32:D2:1537:G:N2	1.98	0.62
1:B0:21:ARG:NH2	1:B0:25:GLU:OE1	2.33	0.62
3:B2:2:LYS:HE2	3:B2:6:LEU:HD13	1.83	0.61
32:D2:774:G:O2'	32:D2:775:G:O5'	2.18	0.61
32:D2:219:A:N3	32:D2:234:U:O2'	2.31	0.61
14:BE:169:LEU:HD22	14:BE:176:PHE:CZ	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:B8:33:THR:OG1	32:D2:2420:C:OP1	2.11	0.61
32:D2:1315:C:O2'	32:D2:1392:A:N3	2.32	0.61
11:BB:216:ARG:NH2	32:D2:781:A:OP1	2.34	0.60
32:D2:1534:U:O2'	32:D2:1537:G:O6	2.05	0.60
15:BF:16:VAL:HG12	15:BF:25:ILE:HG12	1.83	0.60
11:BB:16:VAL:HB	11:BB:203:VAL:HG22	1.84	0.60
27:BU:31:GLU:N	27:BU:31:GLU:OE1	2.34	0.60
32:D2:240:C:OP2	32:D2:241:A:O2'	2.10	0.60
25:BS:32:VAL:HG22	25:BS:37:LYS:HG3	1.84	0.60
1:B0:39:THR:HG21	32:D2:2336:A:H61	1.67	0.60
17:BI:72:ILE:HB	17:BI:108:VAL:HG22	1.83	0.60
32:D2:227:A:O2'	32:D2:228:C:O4'	2.20	0.60
17:BI:7:ASP:OD1	17:BI:8:LYS:N	2.35	0.60
18:BK:100:ILE:O	18:BK:140:GLU:N	2.35	0.60
1:B0:39:THR:HG22	32:D2:2331:G:O2'	2.02	0.59
2:B1:64:ASP:OD1	2:B1:65:THR:N	2.33	0.59
15:BF:37:ASN:OD1	15:BF:38:ASP:N	2.36	0.59
32:D2:1093:G:H21	32:D2:1098:A:H62	1.48	0.59
32:D2:2299:U:O4	32:D2:2318:G:N2	2.36	0.59
33:D3:30:C:H1'	33:D3:57:A:H61	1.68	0.59
15:BF:1:SER:OG	15:BF:2:ARG:N	2.27	0.59
14:BE:89:THR:OG1	33:D3:42:C:N3	2.35	0.59
32:D2:2111:U:O5'	32:D2:2118:U:O2'	2.20	0.58
32:D2:33:C:N4	32:D2:446:G:O2'	2.36	0.58
32:D2:1083:U:O2'	32:D2:1085:A:N7	2.26	0.58
21:BO:41:ARG:NH2	32:D2:807:U:OP2	2.36	0.58
24:BR:56:LYS:O	24:BR:59:ALA:N	2.36	0.58
32:D2:2291:U:OP1	32:D2:2380:C:O2'	2.20	0.58
14:BE:3:LEU:HD12	14:BE:100:GLU:HG3	1.84	0.58
15:BF:88:LEU:O	15:BF:128:THR:OG1	2.22	0.58
33:D3:24:G:N7	33:D3:56:G:O2'	2.27	0.58
24:BR:34:HIS:ND1	24:BR:53:THR:OG1	2.33	0.57
32:D2:2788:C:O2'	32:D2:2809:A:N3	2.34	0.57
17:BI:8:LYS:NZ	17:BI:137:GLU:OE1	2.37	0.57
32:D2:1076:C:O2'	32:D2:1077:A:O4'	2.22	0.57
23:BQ:114:GLU:OE1	23:BQ:118:ARG:NH1	2.37	0.57
10:B9:36:ARG:HG2	10:B9:37:GLN:H	1.69	0.57
27:BU:85:LYS:NZ	32:D2:815:C:OP1	2.37	0.57
32:D2:1070:A:N7	32:D2:1096:A:O2'	2.38	0.57
32:D2:1730:C:O2	32:D2:1731:G:N1	2.38	0.57
1:B0:29:ALA:N	1:B0:60:ASP:OD1	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:BK:75:ALA:N	32:D2:1059:G:O2'	2.37	0.57
22:BP:106:ASP:OD2	22:BP:107:GLY:N	2.38	0.57
32:D2:631:A:N3	32:D2:2415:G:O2'	2.35	0.57
34:FS:17:LEU:O	34:FS:47:HIS:NE2	2.37	0.57
16:BG:35:VAL:HG22	16:BG:50:ALA:HB1	1.87	0.56
17:BI:15:LEU:O	17:BI:15:LEU:HD23	2.06	0.56
31:BY:76:ASP:OD1	31:BY:77:VAL:N	2.39	0.56
32:D2:2135:A:N6	32:D2:2156:G:O2'	2.38	0.56
3:B2:18:LEU:HD11	29:BW:4:GLU:OE1	2.05	0.56
32:D2:2508:G:HO2'	32:D2:2554:U:HO2'	1.54	0.56
32:D2:1172:C:O2'	32:D2:1173:U:O4'	2.23	0.56
16:BG:35:VAL:HG21	16:BG:54:MET:HG3	1.87	0.56
29:BW:56:GLU:HG3	29:BW:57:VAL:HG23	1.88	0.56
32:D2:2031:A:N3	32:D2:2455:G:O2'	2.37	0.55
11:BB:143:VAL:HG12	11:BB:189:ALA:CB	2.36	0.55
22:BP:44:ARG:NH1	32:D2:2484:G:OP1	2.36	0.55
32:D2:1668:A:O2'	32:D2:1674:G:N7	2.34	0.55
32:D2:1833:C:O2'	32:D2:1969:A:N1	2.38	0.55
32:D2:1969:A:O2'	32:D2:1972:G:N3	2.30	0.55
26:BT:103:VAL:O	26:BT:106:THR:OG1	2.17	0.55
11:BB:177:SER:OG	32:D2:1799:G:N7	2.39	0.55
16:BG:60:ARG:NE	32:D2:341:C:OP1	2.39	0.55
32:D2:2655:G:N2	32:D2:2665:A:OP2	2.39	0.55
13:BD:87:ALA:O	13:BD:88:ARG:NH1	2.40	0.55
14:BE:37:MET:HE3	14:BE:52:ALA:HB1	1.89	0.55
14:BE:128:SER:HG	32:D2:2303:G:HO2'	1.54	0.55
30:BX:80:ASP:OD1	30:BX:81:ARG:N	2.38	0.55
18:BK:123:ALA:HB1	32:D2:1081:U:H4'	1.88	0.55
27:BU:14:VAL:HG12	27:BU:20:VAL:HG11	1.89	0.55
32:D2:1450:G:H21	32:D2:1452:G:H1	1.55	0.55
32:D2:1952:A:N3	32:D2:2560:A:O2'	2.37	0.54
34:FS:11:ILE:O	34:FS:14:ILE:HG22	2.07	0.54
9:B8:4:LYS:NZ	32:D2:254:G:N7	2.55	0.54
12:BC:183:GLU:OE1	12:BC:183:GLU:N	2.38	0.54
25:BS:1:SER:OG	25:BS:2:ASN:N	2.33	0.54
32:D2:1528:A:OP2	32:D2:1543:G:N2	2.39	0.54
32:D2:1964:G:O2'	32:D2:1967:C:OP2	2.21	0.54
13:BD:53:THR:HG21	32:D2:452:G:C8	2.42	0.54
27:BU:55:ASP:OD2	27:BU:56:GLY:N	2.37	0.54
32:D2:2144:G:H1'	32:D2:2147:A:H61	1.73	0.54
8:B7:24:THR:HG23	8:B7:27:GLY:H	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:BD:111:GLU:OE2	13:BD:115:GLN:NE2	2.40	0.54
14:BE:56:LEU:HD22	14:BE:64:PRO:HB3	1.89	0.54
32:D2:2134:A:C5	32:D2:2156:G:N2	2.76	0.53
1:B0:67:VAL:HG12	1:B0:67:VAL:O	2.09	0.53
25:BS:26:GLU:OE2	25:BS:41:ALA:HB1	2.09	0.53
14:BE:3:LEU:HD11	14:BE:103:ILE:HD11	1.90	0.53
4:B3:23:LEU:HD11	4:B3:53:MET:SD	2.48	0.53
6:B5:37:HIS:ND1	6:B5:38:LEU:O	2.42	0.53
32:D2:1125:G:OP2	32:D2:1126:A:O2'	2.18	0.53
32:D2:1417:C:O2'	32:D2:1587:G:O2'	2.18	0.53
11:BB:1:ALA:N	11:BB:19:VAL:O	2.38	0.53
13:BD:53:THR:HG21	32:D2:452:G:H8	1.74	0.53
4:B3:44:ARG:NH2	4:B3:58:GLU:OE2	2.40	0.52
13:BD:113:VAL:HG22	13:BD:118:LEU:HD23	1.90	0.52
28:BV:65:ASP:OD1	28:BV:68:ASP:N	2.38	0.52
29:BW:29:THR:OG1	29:BW:86:THR:HG22	2.08	0.52
20:BN:38:ILE:HD13	20:BN:61:VAL:HG22	1.90	0.52
34:FS:14:ILE:HG21	34:FS:22:ILE:HG12	1.91	0.52
26:BT:5:ARG:NH2	32:D2:585:G:N7	2.58	0.52
19:BM:37:ARG:NH1	32:D2:1007:C:OP1	2.43	0.52
32:D2:1173:U:O2	32:D2:1177:G:N1	2.40	0.52
32:D2:141:G:OP2	32:D2:141:G:N2	2.29	0.52
4:B3:6:ILE:CD1	4:B3:47:ILE:HD11	2.39	0.52
10:B9:16:ILE:HD13	10:B9:25:VAL:HG22	1.92	0.51
14:BE:7:TYR:HA	14:BE:11:VAL:HG22	1.92	0.51
15:BF:24:THR:HG22	15:BF:31:GLU:OE2	2.11	0.51
17:BI:115:VAL:HG12	17:BI:132:PHE:CE2	2.45	0.51
32:D2:1063:G:O6	32:D2:1075:C:N4	2.43	0.51
32:D2:1378:A:O2'	32:D2:1379:U:OP2	2.23	0.51
32:D2:1966:A:N3	32:D2:2592:G:O2'	2.37	0.51
5:B4:28:VAL:HG21	5:B4:32:LEU:HD11	1.92	0.51
17:BI:143:ILE:HG23	17:BI:143:ILE:O	2.09	0.51
32:D2:2857:G:N2	32:D2:2860:A:OP2	2.33	0.51
22:BP:123:LYS:NZ	32:D2:2483:C:N3	2.58	0.51
31:BY:43:ASP:OD1	31:BY:44:HIS:N	2.44	0.51
32:D2:1454:C:O2'	32:D2:1455:G:O5'	2.23	0.51
15:BF:51:PHE:HE1	15:BF:71:LEU:HD12	1.75	0.51
32:D2:500:G:N1	32:D2:503:A:OP2	2.41	0.51
2:B1:2:ARG:NH1	32:D2:1365:A:OP1	2.41	0.51
18:BK:127:SER:OG	32:D2:1059:G:N2	2.44	0.50
32:D2:463:G:N2	32:D2:466:A:OP2	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:D2:1838:C:N4	32:D2:1899:A:OP2	2.38	0.50
11:BB:70:LYS:NZ	11:BB:97:ASP:OD2	2.41	0.50
13:BD:149:ILE:HD11	13:BD:188:MET:SD	2.52	0.50
32:D2:2128:G:H21	32:D2:2173:A:H1'	1.76	0.50
21:BO:95:LEU:HD22	21:BO:100:ILE:HD11	1.94	0.50
34:FS:40:CYS:HB3	34:FS:89:VAL:HG12	1.92	0.50
18:BK:133:ARG:NH2	32:D2:1079:C:O2'	2.45	0.50
9:B8:38:LYS:NZ	32:D2:2365:G:N7	2.59	0.50
12:BC:3:GLY:C	12:BC:4:LEU:HD12	2.37	0.50
17:BI:70:GLU:OE1	17:BI:70:GLU:N	2.40	0.50
5:B4:15:SER:O	5:B4:34:LEU:N	2.45	0.50
15:BF:48:THR:OG1	15:BF:49:LEU:N	2.45	0.50
1:B0:38:GLY:HA2	32:D2:2330:G:H21	1.77	0.50
34:FS:42:GLY:N	34:FS:90:MET:O	2.42	0.50
12:BC:151:THR:OG1	32:D2:2032:G:N2	2.45	0.49
32:D2:2144:G:O2'	32:D2:2147:A:N1	2.42	0.49
32:D2:1081:U:N3	32:D2:1082:U:O4	2.44	0.49
32:D2:1019:U:OP1	32:D2:1035:U:O2'	2.23	0.49
33:D3:88:C:HO2'	33:D3:89:U:P	2.31	0.49
32:D2:1093:G:N2	32:D2:1098:A:H62	2.09	0.49
28:BV:42:LYS:O	28:BV:45:VAL:HG12	2.13	0.49
32:D2:774:G:HO2'	32:D2:775:G:P	2.35	0.49
24:BR:111:ARG:NH2	32:D2:2376:A:N3	2.61	0.49
32:D2:837:C:N3	32:D2:941:A:N6	2.60	0.49
32:D2:1028:A:N3	32:D2:2486:C:O2'	2.38	0.49
13:BD:149:ILE:HG22	13:BD:170:ARG:HB2	1.94	0.49
16:BG:71:LYS:NZ	32:D2:613:A:OP2	2.43	0.49
14:BE:70:ARG:NH1	32:D2:2298:A:OP1	2.44	0.48
23:BQ:53:THR:HG21	32:D2:2840:C:C5'	2.41	0.48
32:D2:886:A:O2'	32:D2:889:C:N4	2.46	0.48
4:B3:4:ILE:HD11	4:B3:56:VAL:HG21	1.95	0.48
32:D2:877:A:O2'	32:D2:900:A:N6	2.46	0.48
18:BK:134:SER:OG	32:D2:1077:A:N3	2.44	0.48
32:D2:966:G:O4'	32:D2:2267:A:N6	2.47	0.48
34:FS:3:GLY:O	34:FS:6:LEU:N	2.46	0.48
32:D2:619:G:OP2	32:D2:620:G:N2	2.43	0.48
28:BV:72:THR:HG21	28:BV:108:SER:OG	2.14	0.48
32:D2:1171:G:O2'	32:D2:1172:C:O5'	2.22	0.48
11:BB:116:GLN:N	11:BB:127:ASN:OD1	2.42	0.48
32:D2:1130:U:N3	32:D2:2025:C:OP1	2.43	0.48
13:BD:176:ASP:OD1	13:BD:176:ASP:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:BE:118:ALA:O	14:BE:166:ARG:NH2	2.46	0.48
3:B2:17:GLU:HB2	3:B2:53:VAL:HG11	1.95	0.48
14:BE:35:LEU:HD13	14:BE:153:ILE:HG12	1.95	0.48
30:BX:67:SER:OG	32:D2:335:C:O2	2.32	0.48
15:BF:104:LEU:HB2	15:BF:112:VAL:HG13	1.95	0.48
24:BR:69:ASP:OD1	24:BR:70:ALA:N	2.47	0.47
32:D2:1385:A:O2'	32:D2:1396:U:O2	2.30	0.47
32:D2:2682:A:H61	32:D2:2728:U:H1'	1.79	0.47
7:B6:39:ASP:O	7:B6:43:ARG:N	2.46	0.47
29:BW:57:VAL:HG12	29:BW:58:VAL:N	2.29	0.47
14:BE:36:ASN:OD1	14:BE:37:MET:N	2.47	0.47
15:BF:8:VAL:O	15:BF:48:THR:OG1	2.23	0.47
32:D2:964:C:O2'	32:D2:2273:A:N3	2.39	0.47
5:B4:28:VAL:HG11	5:B4:32:LEU:HD21	1.97	0.47
19:BM:49:ASP:OD1	19:BM:121:LYS:NZ	2.34	0.47
32:D2:2075:U:OP2	32:D2:2238:G:O2'	2.32	0.47
1:B0:36:GLN:OE1	1:B0:40:LYS:N	2.47	0.47
12:BC:114:LYS:NZ	32:D2:2723:C:OP1	2.45	0.47
22:BP:53:MET:HE1	22:BP:103:TYR:CD1	2.50	0.47
33:D3:3:C:O2'	33:D3:4:C:OP1	2.30	0.47
2:B1:60:LYS:NZ	32:D2:371:A:O2'	2.43	0.47
14:BE:140:ILE:HD12	14:BE:145:VAL:HG21	1.96	0.47
21:BO:77:ILE:HD11	21:BO:110:VAL:HG22	1.97	0.47
25:BS:96:LEU:HB3	25:BS:99:LEU:HD23	1.96	0.47
28:BV:7:HIS:CD2	28:BV:10:ALA:HB2	2.50	0.47
28:BV:61:ASN:ND2	32:D2:495:G:N3	2.59	0.47
30:BX:97:SER:O	30:BX:98:ASN:OD1	2.32	0.47
32:D2:1536:C:O2'	32:D2:1537:G:N2	2.47	0.47
32:D2:1715:G:O2'	32:D2:1716:U:O5'	2.32	0.47
17:BI:122:LEU:HD23	17:BI:122:LEU:O	2.14	0.47
18:BK:18:ASN:ND2	18:BK:38:CYS:SG	2.88	0.47
32:D2:586:A:N1	32:D2:809:G:O2'	2.40	0.47
28:BV:4:ILE:HG12	28:BV:106:VAL:HG22	1.97	0.47
32:D2:1171:G:N1	32:D2:1179:G:O6	2.48	0.47
11:BB:250:GLN:NE2	11:BB:251:THR:O	2.48	0.46
32:D2:1667:G:O2'	32:D2:1991:U:O4	2.32	0.46
14:BE:65:LEU:HD22	33:D3:42:C:C6	2.50	0.46
32:D2:878:A:N6	32:D2:899:A:O2'	2.48	0.46
2:B1:73:ARG:NH2	2:B1:75:GLU:OE2	2.48	0.46
20:BN:43:ILE:HD12	20:BN:56:ASP:HB2	1.98	0.46
27:BU:49:ILE:O	27:BU:49:ILE:HG13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B6:36:LYS:HG2	7:B6:47:ILE:HD13	1.97	0.46
12:BC:156:PHE:CD1	19:BM:81:ILE:HD13	2.51	0.46
14:BE:139:GLU:OE1	14:BE:139:GLU:N	2.48	0.46
13:BD:127:GLU:O	13:BD:156:ASN:ND2	2.49	0.46
7:B6:36:LYS:CG	7:B6:47:ILE:HD13	2.45	0.46
32:D2:1826:G:O2'	32:D2:1971:U:OP2	2.34	0.46
1:B0:51:ARG:NH1	32:D2:2364:C:OP1	2.48	0.46
2:B1:61:LYS:NZ	32:D2:372:G:OP2	2.42	0.46
11:BB:259:ASN:C	11:BB:259:ASN:HD22	2.24	0.46
10:B9:2:LYS:NZ	10:B9:32:LYS:O	2.45	0.46
11:BB:129:LEU:HD11	11:BB:134:ILE:HG13	1.98	0.46
12:BC:124:ARG:HA	12:BC:165:MET:HE3	1.98	0.46
23:BQ:23:ASN:ND2	32:D2:1294:U:O2	2.49	0.45
16:BG:19:LYS:HE2	16:BG:66:LEU:HD21	1.97	0.45
20:BN:86:LEU:C	20:BN:87:LEU:HD22	2.40	0.45
24:BR:63:LYS:NZ	33:D3:51:G:OP1	2.35	0.45
32:D2:1730:C:O2'	32:D2:1731:G:O5'	2.33	0.45
11:BB:44:ASN:ND2	32:D2:1370:C:O3'	2.44	0.45
27:BU:61:ALA:HB1	27:BU:96:VAL:HG22	1.97	0.45
11:BB:145:MET:CE	11:BB:153:LEU:HD11	2.44	0.45
20:BN:87:LEU:HD12	20:BN:92:GLU:O	2.16	0.45
32:D2:1416:G:O2'	32:D2:1417:C:O5'	2.30	0.45
20:BN:105:ARG:HH12	20:BN:122:VAL:HG12	1.81	0.45
22:BP:23:GLY:O	22:BP:101:VAL:HG22	2.17	0.45
32:D2:247:G:OP2	32:D2:249:C:N4	2.44	0.45
34:FS:14:ILE:HD13	34:FS:22:ILE:HG23	1.99	0.45
34:FS:38:ILE:HB	34:FS:87:VAL:HG22	1.98	0.45
33:D3:66:A:OP2	33:D3:108:A:N6	2.49	0.45
15:BF:38:ASP:OD1	15:BF:39:ALA:N	2.50	0.45
32:D2:987:C:O2'	32:D2:1000:A:N3	2.47	0.45
19:BM:39:LYS:NZ	32:D2:1009:A:OP1	2.50	0.45
25:BS:102:ARG:NH2	32:D2:1754:A:O3'	2.46	0.45
5:B4:1:MET:N	33:D3:44:G:OP2	2.42	0.44
12:BC:83:ARG:NH1	32:D2:2637:U:OP1	2.50	0.44
14:BE:162:ASP:N	14:BE:162:ASP:OD1	2.50	0.44
15:BF:31:GLU:C	15:BF:32:LEU:HD22	2.42	0.44
18:BK:105:LEU:HD23	18:BK:108:ILE:HD12	1.99	0.44
24:BR:79:ALA:O	24:BR:83:LEU:HD23	2.17	0.44
32:D2:177:G:OP2	32:D2:177:G:N2	2.38	0.44
32:D2:1240:U:O2'	32:D2:1241:A:O5'	2.34	0.44
10:B9:7:VAL:HG22	10:B9:38:GLY:HA3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:BP:12:MET:O	22:BP:86:LYS:NZ	2.51	0.44
12:BC:136:ASN:ND2	12:BC:139:SER:O	2.50	0.44
23:BQ:2:ARG:NH2	32:D2:2820:A:OP2	2.50	0.44
13:BD:41:GLN:NE2	32:D2:442:G:O4'	2.50	0.44
28:BV:1:MET:SD	28:BV:1:MET:N	2.83	0.44
32:D2:421:C:HO2'	32:D2:422:A:P	2.41	0.44
27:BU:43:ASN:OD1	27:BU:44:GLY:N	2.45	0.44
11:BB:60:ALA:O	11:BB:62:ARG:NH1	2.51	0.44
32:D2:1418:G:N2	32:D2:1579:A:N7	2.66	0.44
14:BE:38:GLY:O	32:D2:2307:G:N2	2.50	0.44
21:BO:126:ARG:NH1	32:D2:635:C:OP2	2.50	0.44
9:B8:63:TYR:HH	32:D2:592:A:HO2'	1.65	0.44
13:BD:164:LEU:HB3	13:BD:167:VAL:HG22	2.00	0.44
18:BK:134:SER:CB	32:D2:1088:A:H61	2.30	0.44
32:D2:394:C:H2'	32:D2:395:U:O4'	2.18	0.44
32:D2:458:G:O2'	32:D2:459:U:OP2	2.36	0.44
17:BI:104:THR:OG1	17:BI:109:GLU:OE1	2.36	0.44
32:D2:1857:G:H22	32:D2:1884:G:H2'	1.82	0.44
18:BK:54:ILE:HG22	18:BK:73:PRO:HG3	2.00	0.43
20:BN:61:VAL:HG21	20:BN:112:PHE:CE1	2.53	0.43
25:BS:95:LYS:NZ	32:D2:2847:U:OP1	2.48	0.43
32:D2:140:C:OP2	32:D2:141:G:N2	2.51	0.43
32:D2:2505:G:O2'	32:D2:2506:U:O5'	2.25	0.43
18:BK:134:SER:OG	32:D2:1088:A:N6	2.51	0.43
32:D2:859:G:O2'	32:D2:860:U:P	2.76	0.43
14:BE:56:LEU:HD23	14:BE:56:LEU:O	2.19	0.43
29:BW:89:GLU:OE1	29:BW:89:GLU:N	2.48	0.43
11:BB:143:VAL:HG12	11:BB:189:ALA:HB2	2.01	0.43
18:BK:20:SER:HB3	18:BK:21:PRO:HD3	2.01	0.43
20:BN:105:ARG:NH2	20:BN:122:VAL:O	2.50	0.43
34:FS:81:ASP:HA	34:FS:86:ILE:HD12	2.01	0.43
24:BR:108:ASP:OD1	24:BR:109:ALA:N	2.51	0.43
32:D2:2848:G:HO2'	32:D2:2867:G:H22	1.61	0.43
11:BB:257:ARG:NH1	32:D2:1799:G:OP1	2.44	0.43
18:BK:87:SER:OG	18:BK:88:GLY:N	2.51	0.43
20:BN:35:VAL:HG12	20:BN:35:VAL:O	2.17	0.43
27:BU:49:ILE:HG22	27:BU:54:VAL:H	1.84	0.43
32:D2:141:G:O2'	32:D2:142:A:OP1	2.28	0.43
4:B3:4:ILE:HD11	4:B3:56:VAL:CG2	2.48	0.43
12:BC:45:TYR:OH	32:D2:2636:C:O2'	2.24	0.42
19:BM:84:ILE:HG23	19:BM:84:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:BO:101:ILE:HG13	21:BO:102:GLY:N	2.34	0.42
28:BV:70:LYS:O	28:BV:107:VAL:HG23	2.19	0.42
17:BI:80:ILE:HD11	17:BI:146:VAL:HB	2.01	0.42
17:BI:80:ILE:HD12	17:BI:94:ILE:CD1	2.49	0.42
27:BU:64:VAL:HG12	27:BU:95:ASP:O	2.19	0.42
19:BM:32:LEU:HD22	19:BM:54:ILE:HG21	2.01	0.42
32:D2:1726:C:H2'	32:D2:1727:C:O4'	2.19	0.42
34:FS:17:LEU:HD22	34:FS:47:HIS:CD2	2.55	0.42
11:BB:65:ASP:OD2	11:BB:101:ARG:NH1	2.49	0.42
22:BP:53:MET:HG3	22:BP:120:ALA:HB2	2.01	0.42
29:BW:47:VAL:CG2	29:BW:55:VAL:HG21	2.49	0.42
32:D2:2323:G:H2'	32:D2:2324:U:O4'	2.19	0.42
32:D2:2405:G:O2'	32:D2:2412:A:N6	2.51	0.42
22:BP:77:PRO:HG2	22:BP:80:VAL:HG21	2.01	0.42
32:D2:577:G:O2'	32:D2:1254:A:OP1	2.38	0.42
3:B2:43:LEU:HA	3:B2:46:VAL:HG12	2.02	0.42
15:BF:165:ASP:OD1	15:BF:165:ASP:N	2.43	0.42
32:D2:647:G:N2	32:D2:2350:C:O2'	2.51	0.42
15:BF:55:ASP:OD2	15:BF:55:ASP:N	2.51	0.42
32:D2:704:G:H2'	32:D2:726:G:H22	1.84	0.42
33:D3:31:C:H1'	33:D3:53:A:H61	1.84	0.42
32:D2:2521:C:O2'	32:D2:2564:A:N3	2.43	0.42
7:B6:22:THR:HG22	7:B6:23:THR:N	2.31	0.42
21:BO:30:THR:HG22	32:D2:810:U:O4	2.20	0.42
29:BW:58:VAL:HG22	29:BW:85:VAL:HG12	2.02	0.42
32:D2:227:A:O2'	32:D2:228:C:P	2.77	0.42
18:BK:9:LYS:NZ	32:D2:1060:U:OP2	2.52	0.41
32:D2:458:G:O2'	32:D2:459:U:P	2.77	0.41
32:D2:1754:A:N1	32:D2:2716:C:O2'	2.48	0.41
32:D2:1907:G:O6	32:D2:1924:C:N4	2.53	0.41
15:BF:48:THR:HG1	15:BF:49:LEU:H	1.64	0.41
25:BS:28:LYS:HE2	25:BS:82:SER:HB2	2.02	0.41
25:BS:113:LEU:HD12	25:BS:113:LEU:O	2.19	0.41
21:BO:125:LEU:H	21:BO:125:LEU:HD23	1.85	0.41
9:B8:11:LYS:NZ	32:D2:249:C:O2	2.53	0.41
23:BQ:52:ILE:HD11	23:BQ:83:LEU:HD13	2.02	0.41
32:D2:1063:G:N2	32:D2:1076:C:O2	2.54	0.41
15:BF:8:VAL:C	15:BF:48:THR:HG1	2.24	0.41
30:BX:24:VAL:HG13	30:BX:33:VAL:CG2	2.51	0.41
32:D2:33:C:O2	32:D2:447:A:N6	2.53	0.41
32:D2:1085:A:O2'	32:D2:1086:A:O4'	2.30	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:D2:2595:G:N2	32:D2:2598:A:OP2	2.46	0.41
15:BF:38:ASP:OD1	15:BF:38:ASP:C	2.63	0.41
22:BP:69:PRO:O	22:BP:70:ASP:OD2	2.38	0.41
20:BN:5:GLN:N	20:BN:21:CYS:O	2.52	0.41
32:D2:2581:G:N1	32:D2:2610:C:O2'	2.52	0.41
4:B3:47:ILE:O	4:B3:51:SER:N	2.54	0.41
9:B8:28:LEU:O	9:B8:28:LEU:HG	2.21	0.41
11:BB:141:HIS:ND1	11:BB:192:GLY:O	2.44	0.41
14:BE:56:LEU:HD21	14:BE:88:VAL:HG12	2.02	0.41
14:BE:143:ASP:OD1	14:BE:143:ASP:N	2.53	0.41
15:BF:173:ALA:O	15:BF:174:LYS:HG2	2.20	0.41
18:BK:112:LYS:HE2	18:BK:116:MET:HE3	2.03	0.41
24:BR:2:ASP:OD1	24:BR:3:LYS:N	2.53	0.41
24:BR:51:ALA:HB3	24:BR:78:VAL:HB	2.03	0.41
29:BW:69:ARG:NH1	29:BW:71:GLY:O	2.54	0.41
32:D2:228:C:N4	32:D2:2407:A:N3	2.69	0.41
13:BD:155:GLU:CB	16:BG:27:MET:HE1	2.51	0.41
32:D2:876:C:H2'	32:D2:877:A:O4'	2.21	0.41
32:D2:1178:C:H2'	32:D2:1179:G:C8	2.56	0.41
32:D2:2884:U:O2'	32:D2:2885:G:O5'	2.38	0.41
6:B5:10:SER:O	6:B5:14:MET:HG3	2.20	0.40
8:B7:9:VAL:HG22	32:D2:1309:G:OP1	2.21	0.40
26:BT:87:VAL:HG12	27:BU:49:ILE:O	2.21	0.40
32:D2:354:A:H2'	32:D2:355:U:O4'	2.21	0.40
32:D2:2162:G:OP1	32:D2:2171:A:N6	2.51	0.40
12:BC:12:THR:OG1	12:BC:13:ARG:N	2.54	0.40
32:D2:2637:U:H2'	32:D2:2638:G:O4'	2.21	0.40
20:BN:64:ARG:NH1	20:BN:102:PRO:O	2.53	0.40
29:BW:61:LEU:HD11	29:BW:82:LYS:HD3	2.03	0.40
32:D2:265:A:N1	32:D2:427:U:O2'	2.52	0.40
34:FS:75:ALA:HA	34:FS:78:ILE:HD11	2.03	0.40
14:BE:110:ILE:HD11	14:BE:113:PHE:CD2	2.56	0.40
22:BP:17:ASN:O	22:BP:38:ARG:NH1	2.55	0.40
18:BK:123:ALA:HB3	18:BK:124:MET:HE2	2.04	0.40
19:BM:30:THR:HG23	32:D2:1006:C:O4'	2.22	0.40
26:BT:85:ALA:O	26:BT:86:SER:OG	2.34	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	
1	B0	73/85 (86%)	68 (93%)	5 (7%)	0	100	100
2	B1	75/78 (96%)	75 (100%)	0	0	100	100
3	B2	61/63 (97%)	59 (97%)	2 (3%)	0	100	100
4	B3	56/59 (95%)	54 (96%)	2 (4%)	0	100	100
5	B4	45/70 (64%)	42 (93%)	3 (7%)	0	100	100
6	B5	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
7	B6	48/55 (87%)	47 (98%)	1 (2%)	0	100	100
8	B7	44/46 (96%)	42 (96%)	2 (4%)	0	100	100
9	B8	62/65 (95%)	58 (94%)	4 (6%)	0	100	100
10	B9	36/38 (95%)	34 (94%)	2 (6%)	0	100	100
11	BB	269/273 (98%)	251 (93%)	18 (7%)	0	100	100
12	BC	207/209 (99%)	198 (96%)	9 (4%)	0	100	100
13	BD	199/201 (99%)	189 (95%)	9 (4%)	1 (0%)	24	48
14	BE	175/179 (98%)	163 (93%)	12 (7%)	0	100	100
15	BF	174/177 (98%)	160 (92%)	14 (8%)	0	100	100
16	BG	69/87 (79%)	67 (97%)	2 (3%)	0	100	100
17	BI	147/149 (99%)	136 (92%)	11 (8%)	0	100	100
18	BK	139/142 (98%)	124 (89%)	15 (11%)	0	100	100
19	BM	140/142 (99%)	139 (99%)	1 (1%)	0	100	100
20	BN	120/123 (98%)	109 (91%)	11 (9%)	0	100	100
21	BO	141/144 (98%)	127 (90%)	14 (10%)	0	100	100
22	BP	134/136 (98%)	129 (96%)	5 (4%)	0	100	100
23	BQ	118/127 (93%)	107 (91%)	11 (9%)	0	100	100
24	BR	114/117 (97%)	109 (96%)	5 (4%)	0	100	100
25	BS	112/115 (97%)	104 (93%)	8 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	BT	115/118 (98%)	112 (97%)	3 (3%)	0	100	100
27	BU	101/103 (98%)	94 (93%)	6 (6%)	1 (1%)	12	32
28	BV	108/110 (98%)	101 (94%)	7 (6%)	0	100	100
29	BW	91/100 (91%)	90 (99%)	1 (1%)	0	100	100
30	BX	100/104 (96%)	89 (89%)	11 (11%)	0	100	100
31	BY	92/94 (98%)	88 (96%)	4 (4%)	0	100	100
34	FS	101/105 (96%)	97 (96%)	4 (4%)	0	100	100
All	All	3520/3671 (96%)	3315 (94%)	203 (6%)	2 (0%)	49	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
27	BU	54	VAL
13	BD	83	VAL

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	
1	B0	57/63 (90%)	57 (100%)	0	100	100
2	B1	67/68 (98%)	67 (100%)	0	100	100
3	B2	55/55 (100%)	55 (100%)	0	100	100
4	B3	48/49 (98%)	48 (100%)	0	100	100
5	B4	43/62 (69%)	43 (100%)	0	100	100
6	B5	47/48 (98%)	47 (100%)	0	100	100
7	B6	45/49 (92%)	45 (100%)	0	100	100
8	B7	38/38 (100%)	38 (100%)	0	100	100
9	B8	51/52 (98%)	51 (100%)	0	100	100
10	B9	34/34 (100%)	34 (100%)	0	100	100
11	BB	216/218 (99%)	216 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
12	BC	164/164 (100%)	164 (100%)	0	100	100
13	BD	165/165 (100%)	165 (100%)	0	100	100
14	BE	148/150 (99%)	148 (100%)	0	100	100
15	BF	137/138 (99%)	137 (100%)	0	100	100
16	BG	54/66 (82%)	54 (100%)	0	100	100
17	BI	114/114 (100%)	114 (100%)	0	100	100
18	BK	109/110 (99%)	109 (100%)	0	100	100
19	BM	116/116 (100%)	116 (100%)	0	100	100
20	BN	103/104 (99%)	103 (100%)	0	100	100
21	BO	102/103 (99%)	102 (100%)	0	100	100
22	BP	109/109 (100%)	109 (100%)	0	100	100
23	BQ	100/103 (97%)	100 (100%)	0	100	100
24	BR	86/87 (99%)	86 (100%)	0	100	100
25	BS	99/100 (99%)	99 (100%)	0	100	100
26	BT	89/90 (99%)	89 (100%)	0	100	100
27	BU	84/84 (100%)	84 (100%)	0	100	100
28	BV	93/93 (100%)	93 (100%)	0	100	100
29	BW	80/84 (95%)	80 (100%)	0	100	100
30	BX	83/85 (98%)	83 (100%)	0	100	100
31	BY	78/78 (100%)	78 (100%)	0	100	100
34	FS	89/91 (98%)	89 (100%)	0	100	100
All	All	2903/2970 (98%)	2903 (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 (17) such sidechains are listed below:

Mol	Chain	Res	Type
2	B1	35	HIS
3	B2	20	ASN
3	B2	38	GLN
4	B3	8	GLN
5	B4	33	ASN
11	BB	89	ASN

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Mol	Chain	Res	Type
13	BD	136	GLN
17	BI	28	ASN
17	BI	33	GLN
20	BN	29	HIS
21	BO	38	GLN
25	BS	55	HIS
26	BT	43	GLN
28	BV	7	HIS
28	BV	9	HIS
31	BY	51	GLN
34	FS	28	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
32	D2	2902/2903 (99%)	465 (16%)	23 (0%)
33	D3	119/120 (99%)	20 (16%)	3 (2%)
All	All	3021/3023 (99%)	485 (16%)	26 (0%)

All (485) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
32	D2	10	A
32	D2	12	U
32	D2	34	U
32	D2	35	G
32	D2	36	G
32	D2	46	G
32	D2	51	G
32	D2	63	A
32	D2	71	A
32	D2	74	A
32	D2	75	G
32	D2	82	U
32	D2	100	U
32	D2	102	U
32	D2	118	A
32	D2	120	U
32	D2	125	A
32	D2	139	U
32	D2	140	C

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Mol	Chain	Res	Type
32	D2	141	G
32	D2	142	A
32	D2	149	A
32	D2	162	U
32	D2	163	C
32	D2	166	U
32	D2	181	A
32	D2	196	A
32	D2	199	A
32	D2	215	G
32	D2	216	A
32	D2	218	A
32	D2	219	A
32	D2	221	A
32	D2	222	A
32	D2	228	C
32	D2	232	G
32	D2	248	G
32	D2	249	C
32	D2	250	G
32	D2	255	A
32	D2	265	A
32	D2	266	G
32	D2	276	U
32	D2	279	A
32	D2	281	C
32	D2	310	A
32	D2	323	C
32	D2	324	A
32	D2	329	G
32	D2	330	A
32	D2	361	G
32	D2	362	A
32	D2	371	A
32	D2	372	G
32	D2	373	U
32	D2	386	G
32	D2	387	U
32	D2	389	G
32	D2	395	U
32	D2	404	A
32	D2	406	G

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Mol	Chain	Res	Type
32	D2	411	G
32	D2	412	A
32	D2	417	C
32	D2	422	A
32	D2	424	G
32	D2	457	A
32	D2	458	G
32	D2	459	U
32	D2	473	G
32	D2	481	G
32	D2	491	G
32	D2	504	A
32	D2	505	A
32	D2	509	C
32	D2	530	G
32	D2	531	C
32	D2	532	A
32	D2	545	U
32	D2	546	U
32	D2	547	A
32	D2	548	G
32	D2	549	G
32	D2	550	C
32	D2	563	A
32	D2	573	U
32	D2	574	A
32	D2	575	A
32	D2	586	A
32	D2	603	A
32	D2	613	A
32	D2	622	G
32	D2	627	A
32	D2	637	A
32	D2	645	C
32	D2	646	U
32	D2	654	A
32	D2	655	A
32	D2	669	G
32	D2	670	A
32	D2	686	U
32	D2	694	U
32	D2	695	G

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Mol	Chain	Res	Type
32	D2	715	A
32	D2	716	A
32	D2	726	G
32	D2	730	A
32	D2	738	G
32	D2	740	C
32	D2	747	C
32	D2	752	A
32	D2	764	A
32	D2	765	C
32	D2	775	G
32	D2	776	G
32	D2	777	G
32	D2	782	A
32	D2	784	G
32	D2	789	A
32	D2	792	A
32	D2	805	G
32	D2	812	C
32	D2	819	A
32	D2	827	U
32	D2	828	U
32	D2	830	G
32	D2	845	A
32	D2	846	U
32	D2	847	U
32	D2	857	G
32	D2	859	G
32	D2	860	U
32	D2	869	G
32	D2	878	A
32	D2	886	A
32	D2	887	U
32	D2	896	A
32	D2	897	C
32	D2	899	A
32	D2	910	A
32	D2	914	G
32	D2	915	C
32	D2	941	A
32	D2	946	C
32	D2	961	C

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Mol	Chain	Res	Type
32	D2	974	G
32	D2	983	A
32	D2	995	C
32	D2	996	A
32	D2	1012	U
32	D2	1013	C
32	D2	1021	A
32	D2	1022	G
32	D2	1026	G
32	D2	1033	U
32	D2	1046	A
32	D2	1060	U
32	D2	1062	G
32	D2	1063	G
32	D2	1065	U
32	D2	1066	U
32	D2	1068	G
32	D2	1069	A
32	D2	1070	A
32	D2	1071	G
32	D2	1072	C
32	D2	1075	C
32	D2	1076	C
32	D2	1078	U
32	D2	1079	C
32	D2	1082	U
32	D2	1083	U
32	D2	1084	A
32	D2	1088	A
32	D2	1090	A
32	D2	1094	U
32	D2	1095	A
32	D2	1103	A
32	D2	1104	C
32	D2	1108	U
32	D2	1110	G
32	D2	1111	A
32	D2	1112	G
32	D2	1131	G
32	D2	1132	U
32	D2	1133	A
32	D2	1135	C

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Mol	Chain	Res	Type
32	D2	1136	G
32	D2	1139	G
32	D2	1142	A
32	D2	1172	C
32	D2	1174	U
32	D2	1175	A
32	D2	1177	G
32	D2	1178	C
32	D2	1179	G
32	D2	1180	U
32	D2	1206	G
32	D2	1212	G
32	D2	1227	G
32	D2	1238	G
32	D2	1250	G
32	D2	1251	C
32	D2	1253	A
32	D2	1256	G
32	D2	1268	A
32	D2	1271	G
32	D2	1272	A
32	D2	1273	U
32	D2	1300	G
32	D2	1301	A
32	D2	1302	A
32	D2	1306	C
32	D2	1329	U
32	D2	1330	C
32	D2	1345	C
32	D2	1352	U
32	D2	1365	A
32	D2	1368	G
32	D2	1379	U
32	D2	1383	A
32	D2	1386	C
32	D2	1395	A
32	D2	1416	G
32	D2	1419	A
32	D2	1420	A
32	D2	1428	C
32	D2	1452	G
32	D2	1454	C

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Mol	Chain	Res	Type
32	D2	1455	G
32	D2	1459	G
32	D2	1460	U
32	D2	1461	C
32	D2	1482	G
32	D2	1490	A
32	D2	1491	G
32	D2	1493	C
32	D2	1508	A
32	D2	1509	A
32	D2	1515	A
32	D2	1524	G
32	D2	1533	C
32	D2	1534	U
32	D2	1535	A
32	D2	1536	C
32	D2	1537	G
32	D2	1540	G
32	D2	1559	U
32	D2	1560	G
32	D2	1569	A
32	D2	1578	U
32	D2	1581	G
32	D2	1583	A
32	D2	1584	U
32	D2	1585	C
32	D2	1608	A
32	D2	1611	C
32	D2	1647	U
32	D2	1648	U
32	D2	1654	A
32	D2	1674	G
32	D2	1677	A
32	D2	1715	G
32	D2	1716	U
32	D2	1729	U
32	D2	1730	C
32	D2	1731	G
32	D2	1732	C
32	D2	1733	G
32	D2	1735	A
32	D2	1738	G

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Mol	Chain	Res	Type
32	D2	1756	G
32	D2	1758	U
32	D2	1764	C
32	D2	1773	A
32	D2	1776	G
32	D2	1780	A
32	D2	1791	A
32	D2	1800	C
32	D2	1801	A
32	D2	1802	A
32	D2	1808	A
32	D2	1816	C
32	D2	1818	U
32	D2	1828	G
32	D2	1829	A
32	D2	1835	G
32	D2	1847	A
32	D2	1857	G
32	D2	1859	U
32	D2	1870	C
32	D2	1871	A
32	D2	1873	G
32	D2	1876	A
32	D2	1884	G
32	D2	1899	A
32	D2	1901	A
32	D2	1906	G
32	D2	1913	A
32	D2	1914	C
32	D2	1919	A
32	D2	1927	A
32	D2	1930	G
32	D2	1938	A
32	D2	1941	C
32	D2	1955	U
32	D2	1963	U
32	D2	1965	C
32	D2	1966	A
32	D2	1967	C
32	D2	1970	A
32	D2	1971	U
32	D2	1972	G

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Mol	Chain	Res	Type
32	D2	1991	U
32	D2	1993	U
32	D2	1997	C
32	D2	2022	U
32	D2	2023	C
32	D2	2027	G
32	D2	2031	A
32	D2	2033	A
32	D2	2043	C
32	D2	2052	A
32	D2	2055	C
32	D2	2056	G
32	D2	2059	A
32	D2	2060	A
32	D2	2061	G
32	D2	2063	C
32	D2	2069	G
32	D2	2093	G
32	D2	2096	C
32	D2	2100	G
32	D2	2101	A
32	D2	2102	G
32	D2	2103	C
32	D2	2104	C
32	D2	2108	A
32	D2	2110	G
32	D2	2111	U
32	D2	2112	G
32	D2	2115	G
32	D2	2118	U
32	D2	2119	A
32	D2	2127	G
32	D2	2128	G
32	D2	2131	U
32	D2	2132	U
32	D2	2133	G
32	D2	2136	G
32	D2	2145	C
32	D2	2157	G
32	D2	2159	G
32	D2	2162	G
32	D2	2164	C

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Mol	Chain	Res	Type
32	D2	2166	U
32	D2	2168	G
32	D2	2170	A
32	D2	2171	A
32	D2	2172	U
32	D2	2173	A
32	D2	2182	U
32	D2	2183	A
32	D2	2185	U
32	D2	2186	G
32	D2	2188	U
32	D2	2192	U
32	D2	2198	A
32	D2	2203	U
32	D2	2204	G
32	D2	2211	A
32	D2	2212	A
32	D2	2213	U
32	D2	2225	A
32	D2	2226	C
32	D2	2238	G
32	D2	2250	G
32	D2	2278	A
32	D2	2283	C
32	D2	2286	G
32	D2	2287	A
32	D2	2303	G
32	D2	2304	G
32	D2	2305	U
32	D2	2307	G
32	D2	2311	A
32	D2	2325	G
32	D2	2327	A
32	D2	2331	G
32	D2	2334	U
32	D2	2335	A
32	D2	2345	G
32	D2	2350	C
32	D2	2361	G
32	D2	2382	G
32	D2	2383	G
32	D2	2385	C

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Mol	Chain	Res	Type
32	D2	2392	A
32	D2	2402	U
32	D2	2406	A
32	D2	2407	A
32	D2	2422	C
32	D2	2424	C
32	D2	2429	G
32	D2	2430	A
32	D2	2435	A
32	D2	2441	U
32	D2	2448	A
32	D2	2476	A
32	D2	2491	U
32	D2	2497	A
32	D2	2498	C
32	D2	2502	G
32	D2	2503	A
32	D2	2505	G
32	D2	2506	U
32	D2	2508	G
32	D2	2518	A
32	D2	2529	G
32	D2	2547	A
32	D2	2552	U
32	D2	2554	U
32	D2	2566	A
32	D2	2567	G
32	D2	2572	A
32	D2	2573	C
32	D2	2576	G
32	D2	2581	G
32	D2	2582	G
32	D2	2602	A
32	D2	2609	U
32	D2	2613	U
32	D2	2615	U
32	D2	2629	U
32	D2	2655	G
32	D2	2682	A
32	D2	2689	U
32	D2	2690	U
32	D2	2714	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
32	D2	2722	G
32	D2	2733	A
32	D2	2744	G
32	D2	2748	A
32	D2	2751	G
32	D2	2764	A
32	D2	2765	A
32	D2	2778	A
32	D2	2779	U
32	D2	2790	U
32	D2	2791	G
32	D2	2797	U
32	D2	2799	A
32	D2	2800	A
32	D2	2809	A
32	D2	2820	A
32	D2	2833	U
32	D2	2834	G
32	D2	2849	U
32	D2	2861	U
32	D2	2867	G
32	D2	2868	A
32	D2	2872	A
32	D2	2880	C
32	D2	2884	U
32	D2	2887	A
33	D3	4	C
33	D3	9	G
33	D3	12	C
33	D3	13	G
33	D3	26	C
33	D3	35	C
33	D3	41	G
33	D3	42	C
33	D3	44	G
33	D3	52	A
33	D3	64	G
33	D3	66	A
33	D3	67	G
33	D3	88	C
33	D3	89	U
33	D3	90	C

Continued on next page...

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Mol	Chain	Res	Type
33	D3	105	G
33	D3	108	A
33	D3	109	A
33	D3	120	A

All (26) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
32	D2	227	A
32	D2	372	G
32	D2	421	C
32	D2	458	G
32	D2	490	C
32	D2	715	A
32	D2	774	G
32	D2	859	G
32	D2	1020	A
32	D2	1130	U
32	D2	1211	C
32	D2	1490	A
32	D2	1715	G
32	D2	1730	C
32	D2	1875	G
32	D2	2109	U
32	D2	2127	G
32	D2	2326	C
32	D2	2333	A
32	D2	2391	G
32	D2	2505	G
32	D2	2798	U
32	D2	2808	G
33	D3	3	C
33	D3	66	A
33	D3	88	C

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

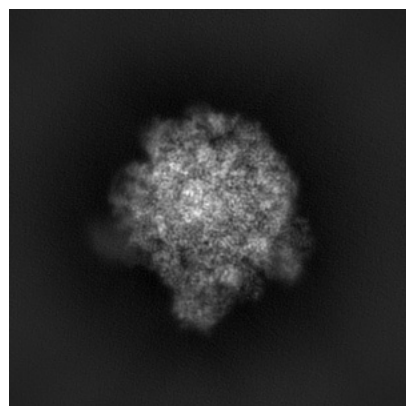
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55177. These allow visual inspection of the internal detail of the map and identification of artifacts.

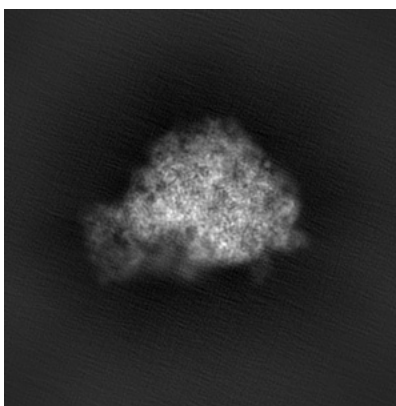
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

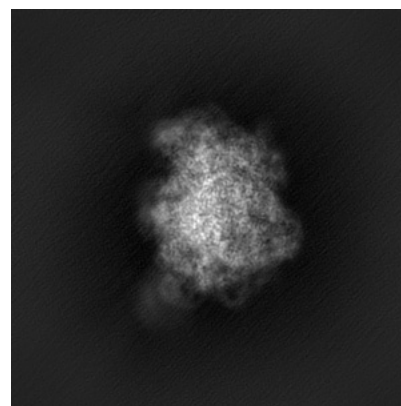
6.1.1 Primary map



X

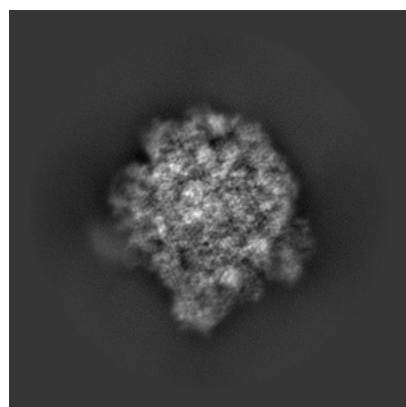


Y

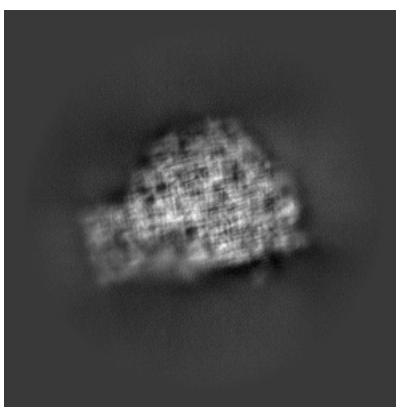


Z

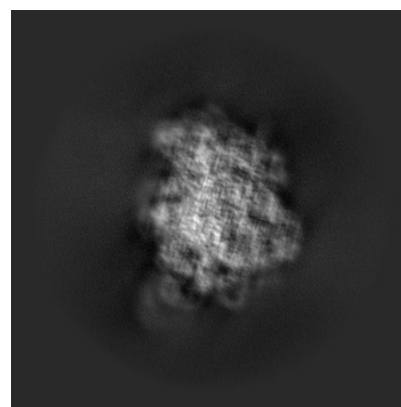
6.1.2 Raw map



X



Y

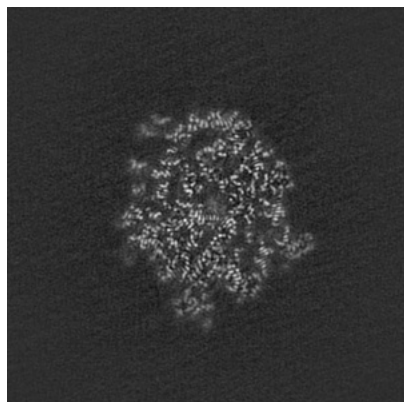


Z

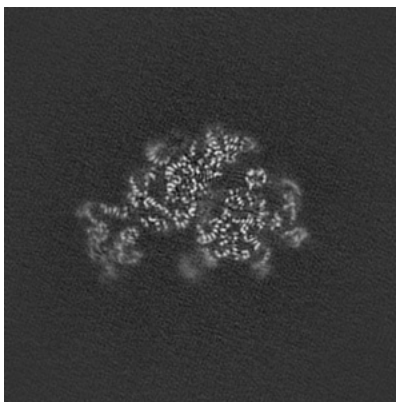
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

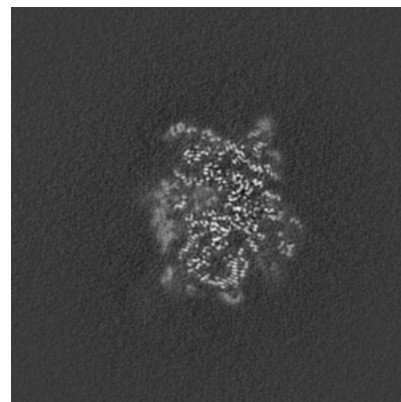
6.2.1 Primary map



X Index: 240

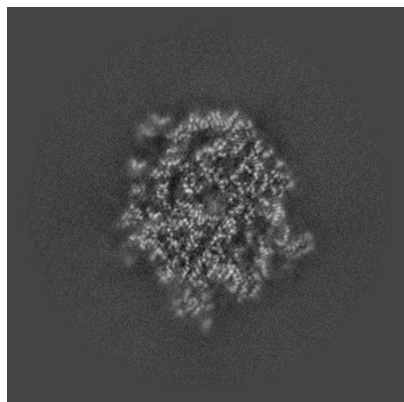


Y Index: 240

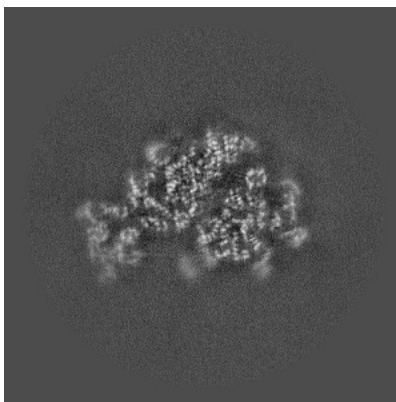


Z Index: 240

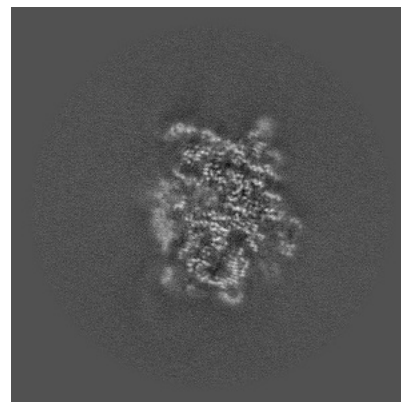
6.2.2 Raw map



X Index: 240



Y Index: 240

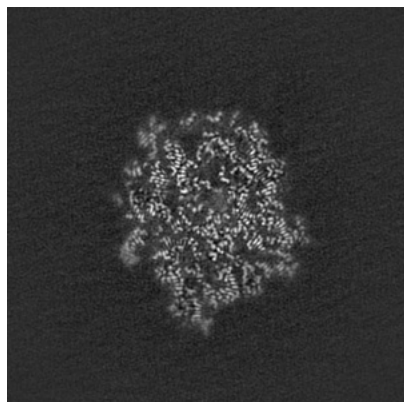


Z Index: 240

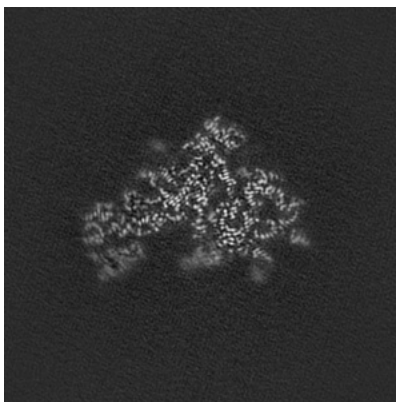
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

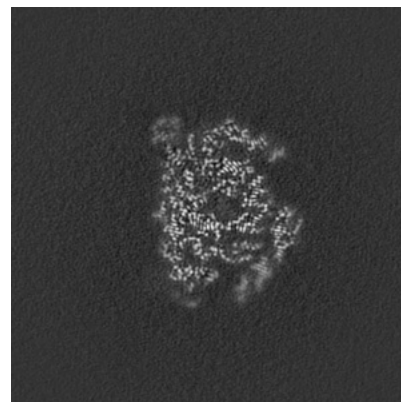
6.3.1 Primary map



X Index: 231

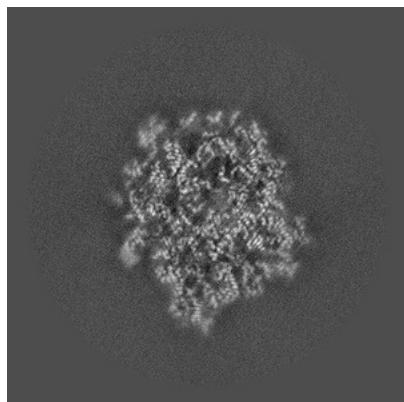


Y Index: 226

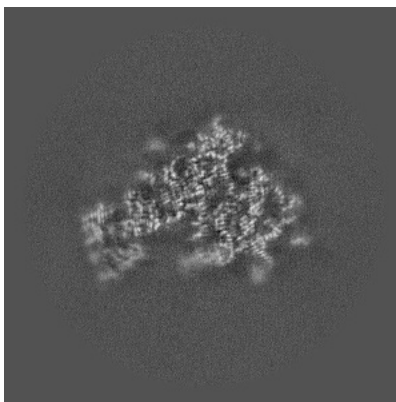


Z Index: 260

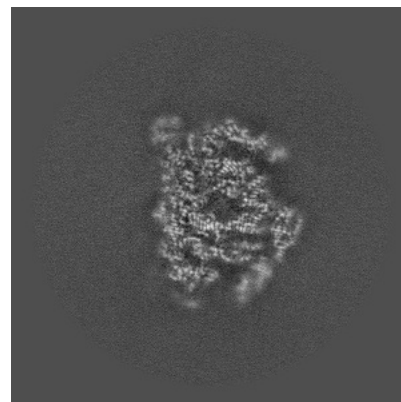
6.3.2 Raw map



X Index: 231



Y Index: 229

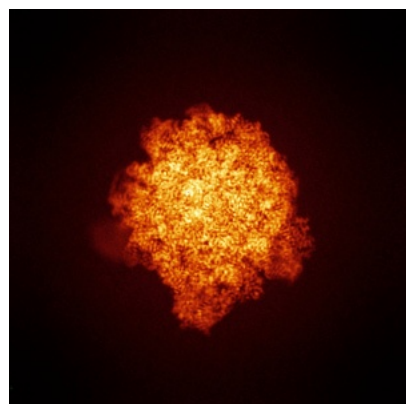


Z Index: 260

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

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

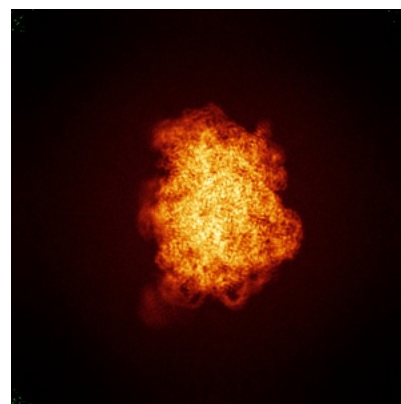
6.4.1 Primary map



X

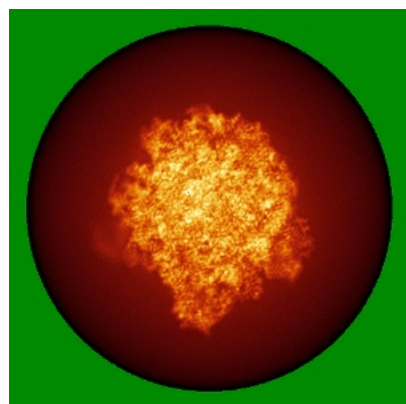


Y

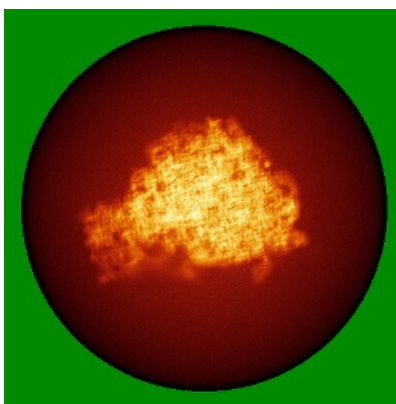


Z

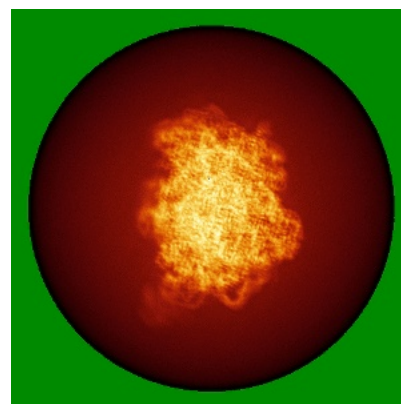
6.4.2 Raw map



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



X



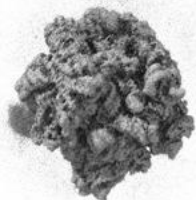
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 2.2. 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



X



Y



Z

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.

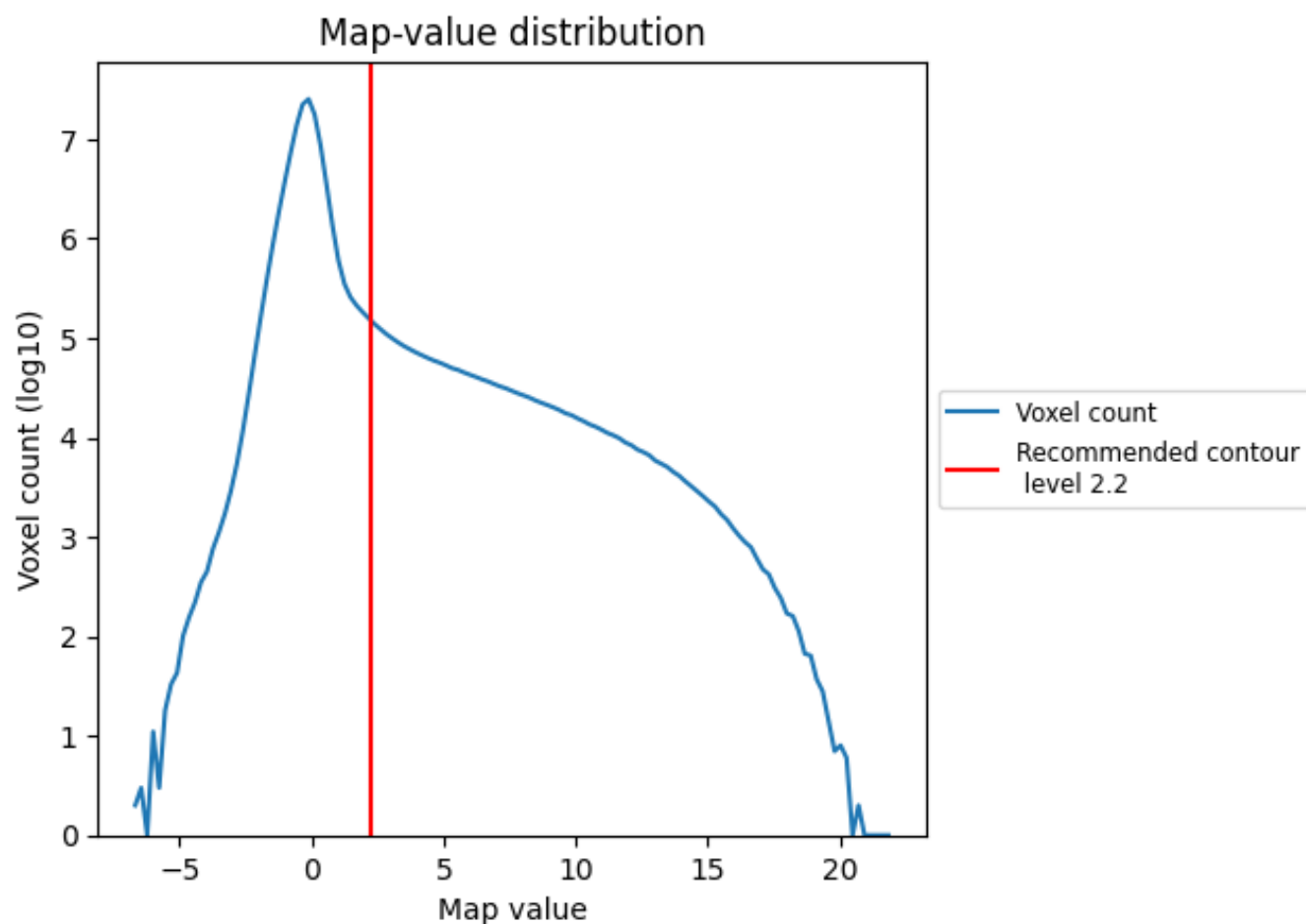
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

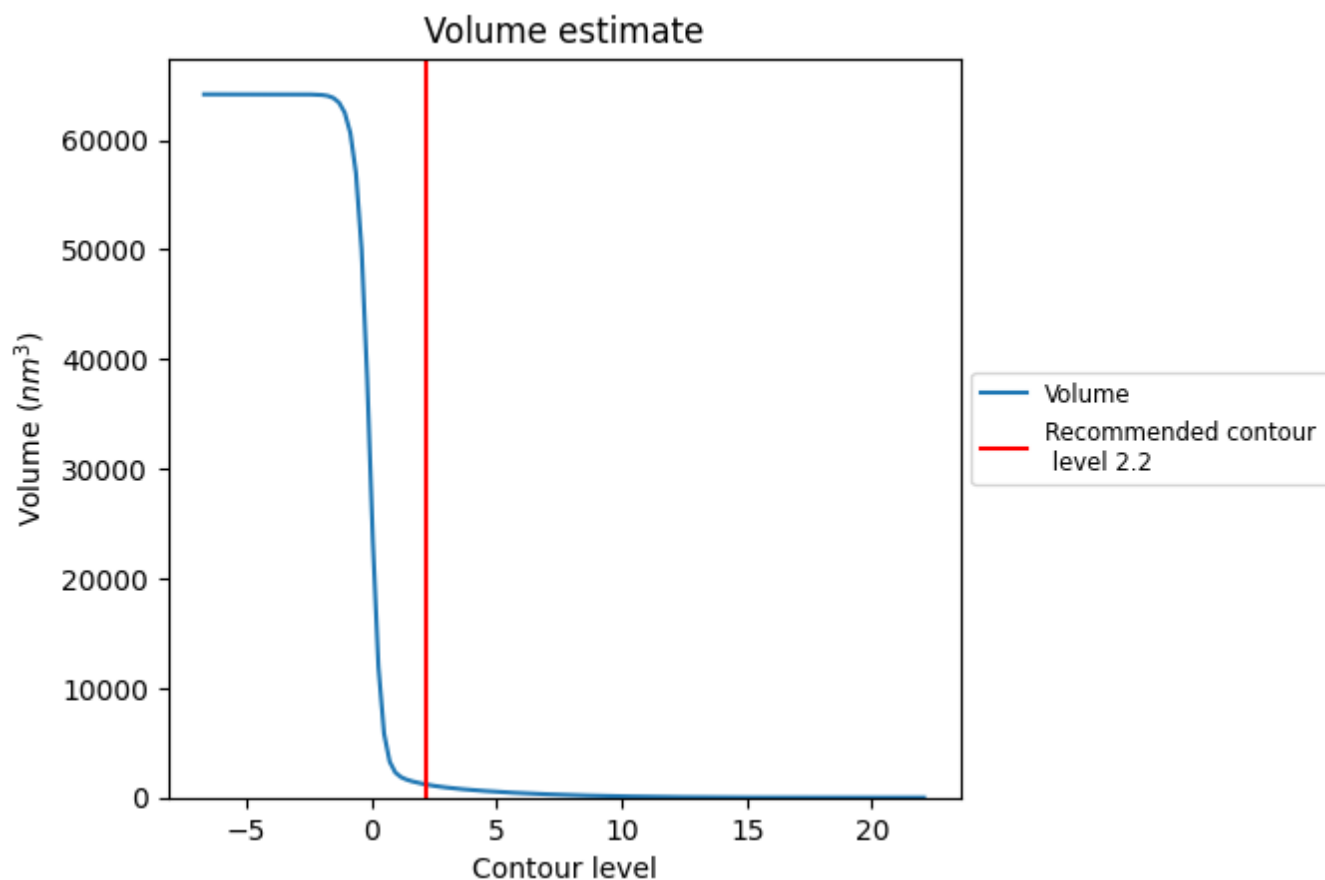
This section contains the results of statistical analysis of the map.

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.

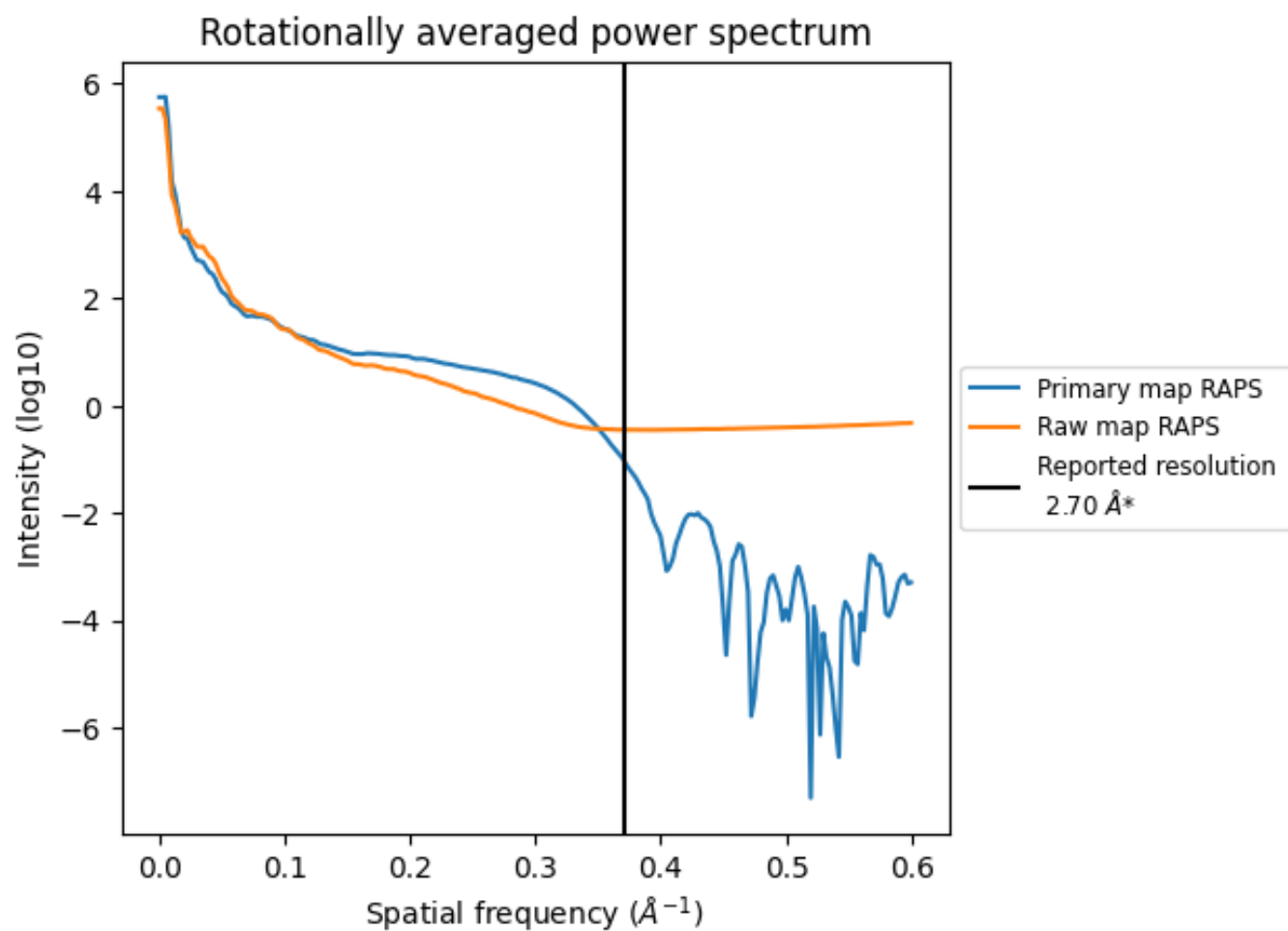
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1188 nm^3 ; this corresponds to an approximate mass of 1074 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 ⓘ

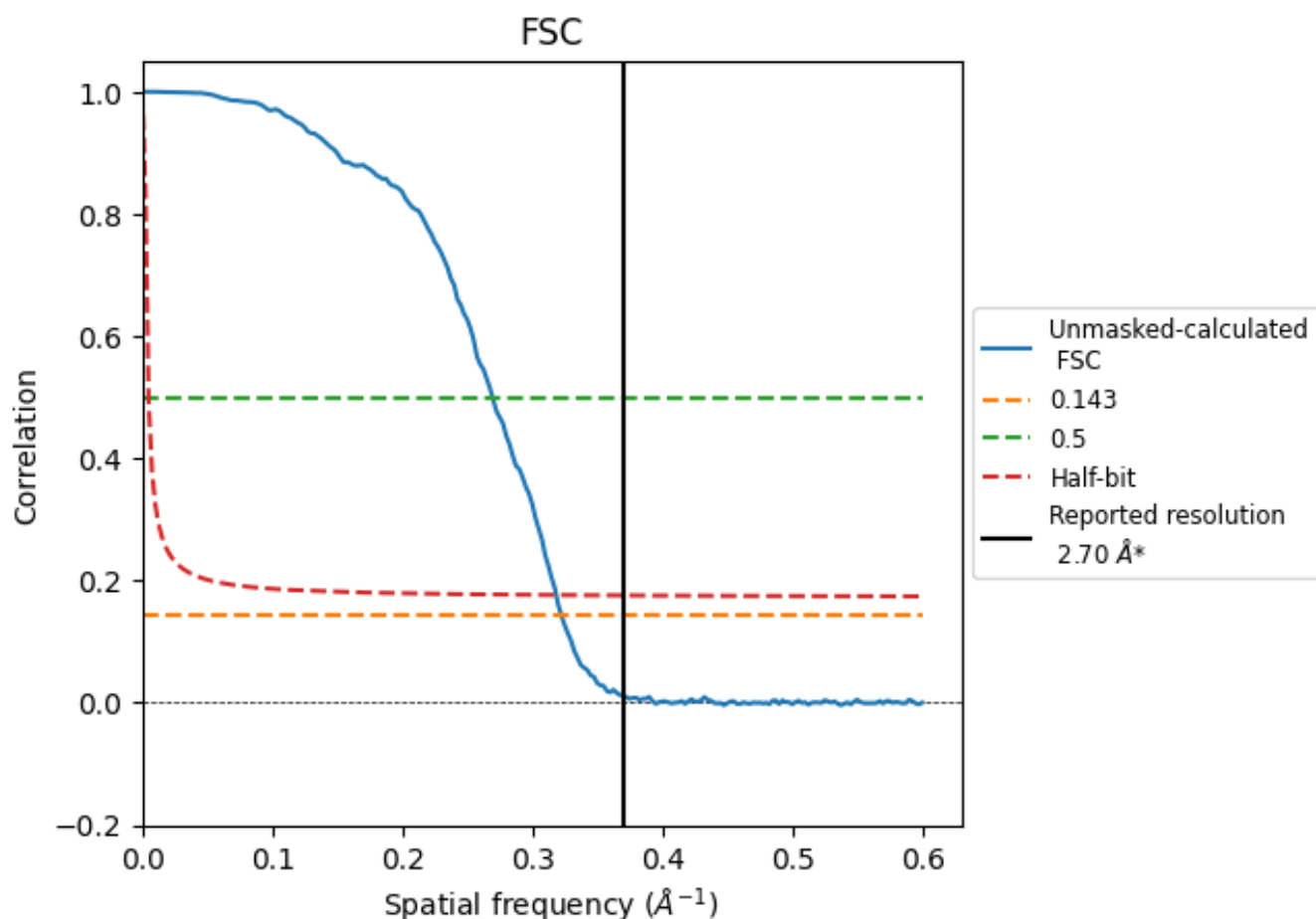


*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

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 0.370 Å⁻¹

8.2 Resolution estimates [i](#)

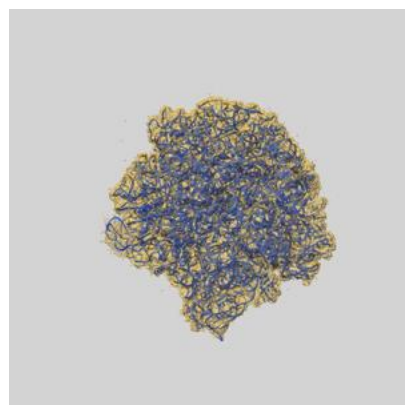
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.10	3.71	3.14

*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 3.10 differs from the reported value 2.7 by more than 10 %

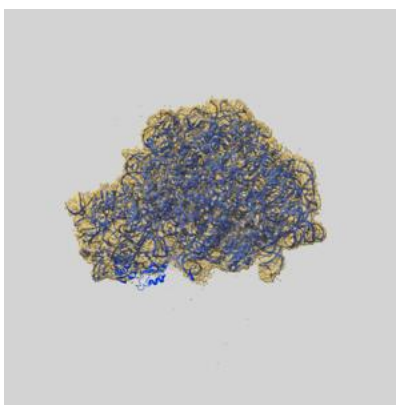
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-55177 and PDB model 9SS5. Per-residue inclusion information can be found in section 3 on page 11.

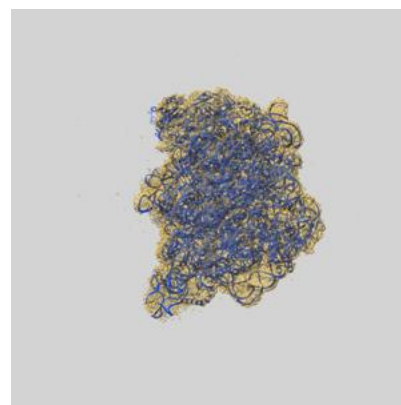
9.1 Map-model overlay [i](#)



X



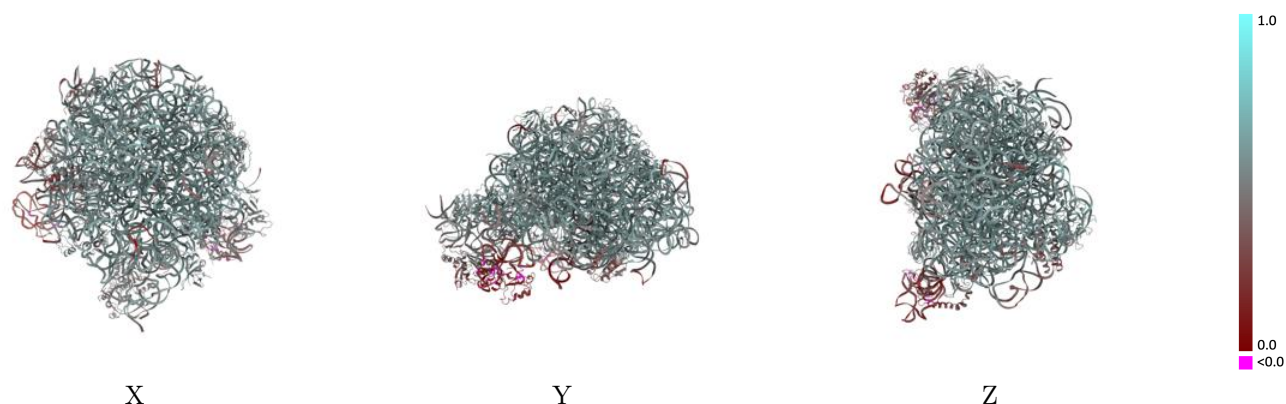
Y



Z

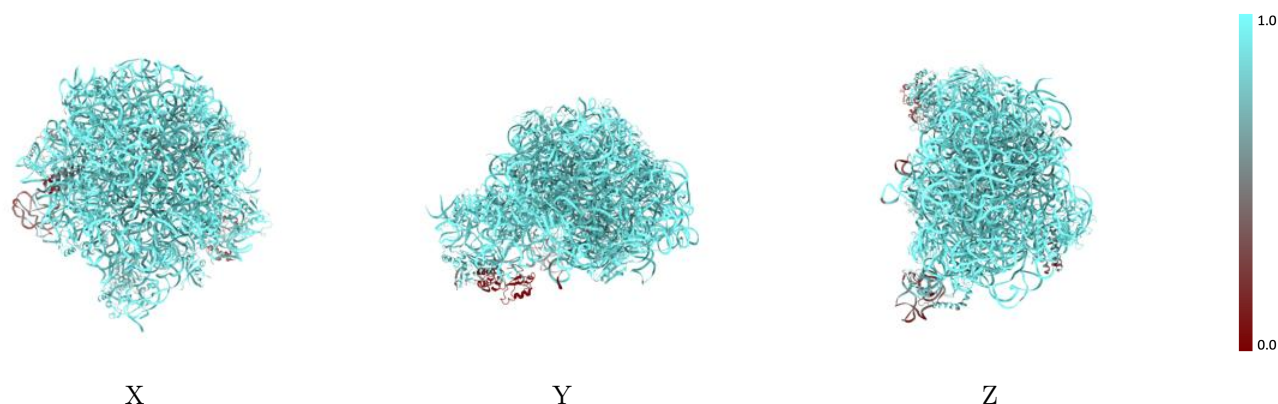
The images above show the 3D surface view of the map at the recommended contour level 2.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



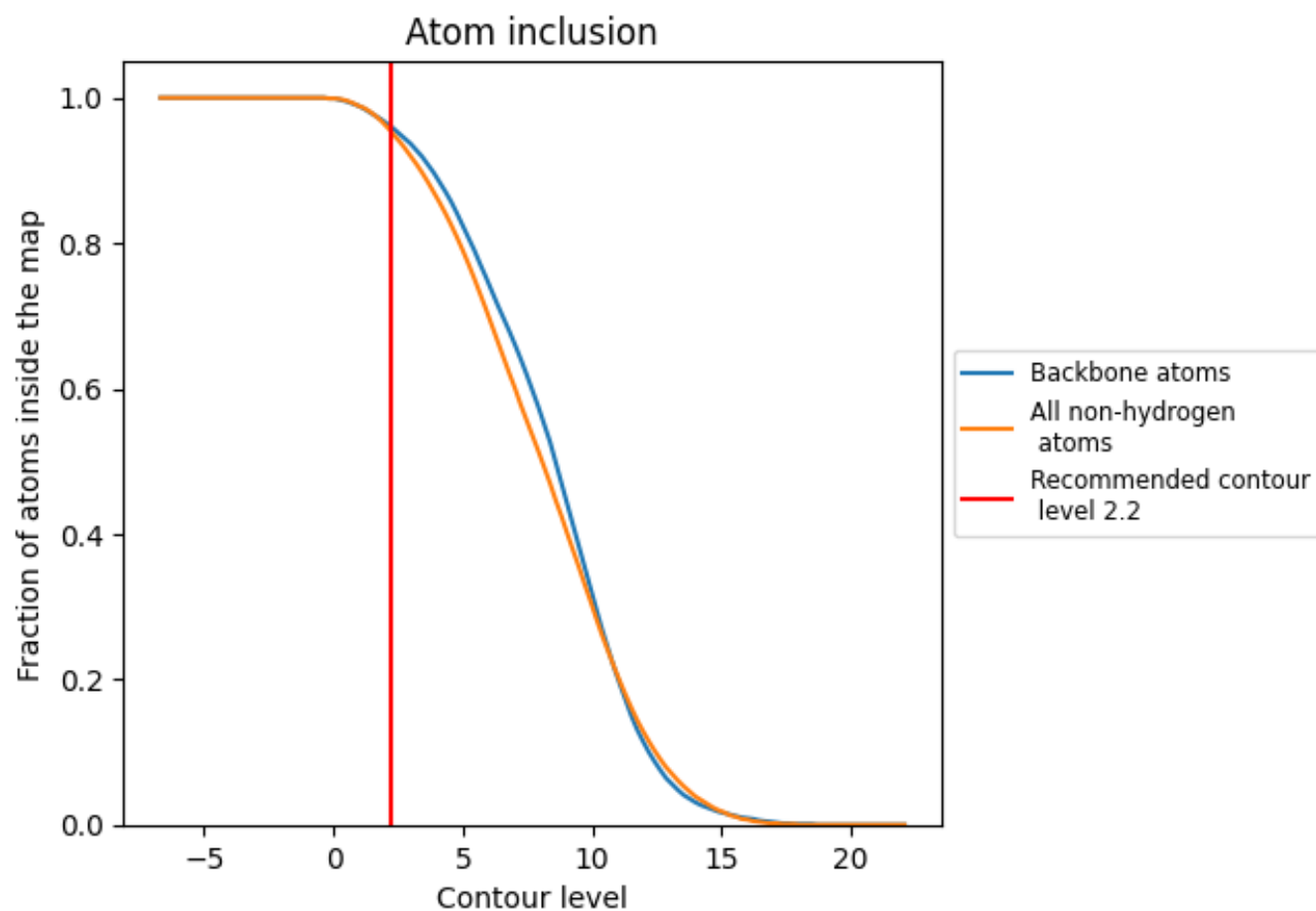
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (2.2).







































































9.4 Atom inclusion ⓘ



At the recommended contour level, 96% of all backbone atoms, 96% 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 (2.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9550	 0.5250
B0	 0.9790	 0.5670
B1	 0.9800	 0.5710
B2	 0.9360	 0.4820
B3	 0.9700	 0.5560
B4	 0.7340	 0.3530
B5	 0.9810	 0.5660
B6	 0.9780	 0.5340
B7	 0.9830	 0.5850
B8	 0.9960	 0.5840
B9	 0.9830	 0.5580
BB	 0.9830	 0.5600
BC	 0.9830	 0.5670
BD	 0.9510	 0.5310
BE	 0.8830	 0.4010
BF	 0.9550	 0.4930
BG	 0.5480	 0.3650
BI	 0.7570	 0.3650
BK	 0.0970	 0.1500
BM	 0.9850	 0.5800
BN	 0.9620	 0.5450
BO	 0.9640	 0.5470
BP	 0.9840	 0.5580
BQ	 0.9840	 0.5670
BR	 0.9490	 0.4880
BS	 0.9550	 0.5280
BT	 0.9890	 0.5780
BU	 0.9560	 0.5460
BV	 0.9680	 0.5640
BW	 0.9640	 0.5470
BX	 0.9840	 0.5400
BY	 0.9550	 0.5360
D2	 0.9720	 0.5330
D3	 0.9930	 0.5160
FS	 0.8750	 0.3800

