



Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 11:41 AM EST

PDB ID : 3J0O
EMDB ID : EMD-5327
Title : Core of mammalian 80S pre-ribosome in complex with tRNAs fitted to a 9A cryo-EM map: classic PRE state 2
Authors : Budkevich, T.; Giesebrecht, J.; Altman, R.; Munro, J.; Mielke, T.; Nierhaus, K.; Blanchard, S.; Spahn, C.M.
Deposited on : 2011-10-05
Resolution : 9.00 Å (reported)
Based on initial models : 2WDK, 3O58, 2XZM

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.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

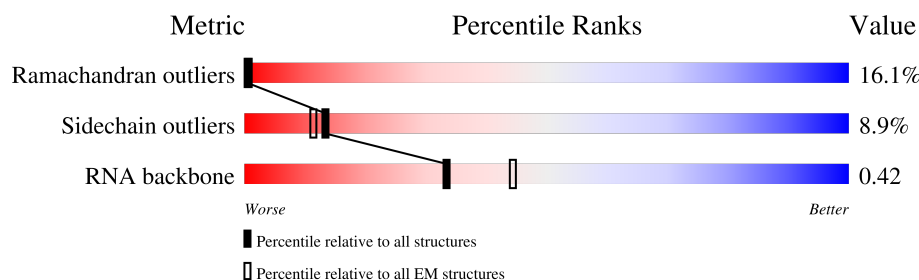
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 9.00 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.
















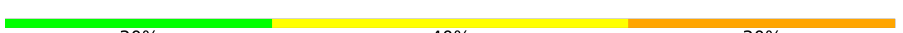








Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	48	<div> <div>65%</div> <div>23%</div> <div>8%</div> <div>.</div> </div>
2	b	12	<div> <div>67%</div> <div>33%</div> </div>
3	c	17	<div> <div>71%</div> <div>29%</div> </div>
4	d	7	<div> <div>71%</div> <div>14%</div> <div>14%</div> </div>
5	e	4	<div> <div>100%</div> </div>
6	E	5	<div> <div>60%</div> <div>20%</div> <div>20%</div> </div>
7	f	21	<div> <div>76%</div> <div>24%</div> </div>
8	g	31	<div> <div>71%</div> <div>23%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
9	G	13	
10	h	111	
11	T	192	
12	K	140	
13	L	141	
14	X	68	
15	S	125	
16	2	112	
17	3	12	
18	4	14	
19	5	6	
20	6	19	
21	7	50	
22	8	20	
23	B	213	
24	F	95	
25	V	76	
25	W	76	
25	Y	76	
26	v	3	
26	y	3	
27	w	2	

2 Entry composition

There are 27 unique types of molecules in this entry. The entry contains 22450 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 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	a	48	Total	C	N	O	P	0	0
			1029	459	190	332	48		

- Molecule 2 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	b	12	Total	C	N	O	P	0	0
			260	116	49	83	12		

- Molecule 3 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	c	17	Total	C	N	O	P	0	0
			362	162	66	117	17		

- Molecule 4 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	d	7	Total	C	N	O	P	0	0
			155	69	33	46	7		

- Molecule 5 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	e	4	Total	C	N	O	P	0	0
			84	38	16	26	4		

- Molecule 6 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	E	5	Total	C	N	O	P	0	0
			100	45	13	37	5		

- Molecule 7 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	f	21	Total	C	N	O	P	0	0
			452	200	79	152	21		

- Molecule 8 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	g	31	Total	C	N	O	P	0	0
			660	295	118	216	31		

- Molecule 9 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	G	13	Total	C	N	O	P	0	0
			276	123	49	91	13		

- Molecule 10 is a RNA chain called 40S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	h	111	Total	C	N	O	P	0	0
			2368	1060	431	766	111		

- Molecule 11 is a protein called Ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	T	192	Total	C	N	O	S	0	0
			1520	961	281	270	8		

- Molecule 12 is a protein called Ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	K	140	Total	C	N	O	S	0	0
			1063	654	206	197	6		

- Molecule 13 is a protein called Ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	L	141	Total	C	N	O	S	0	0
			1097	691	221	180	5		

- Molecule 14 is a protein called Ribosomal protein S30.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	X	68	Total	C	N	O	S	0	0
			554	350	113	90	1		

- Molecule 15 is a protein called Ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	S	125	Total	C	N	O	S	0	0
			985	632	173	176	4		

- Molecule 16 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	2	112	Total	C	N	O	P	0	0
			2392	1070	435	775	112		

- Molecule 17 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	3	12	Total	C	N	O	P	0	0
			259	116	50	81	12		

- Molecule 18 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	4	14	Total	C	N	O	P	0	0
			306	135	59	98	14		

- Molecule 19 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	5	6	Total	C	N	O	P	0	0
			127	57	23	41	6		

- Molecule 20 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	6	19	Total	C	N	O	P	0	0
			417	187	88	123	19		

- Molecule 21 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	7	50	Total	C	N	O	P	0	0
			1054	471	173	360	50		

- Molecule 22 is a RNA chain called 60S ribosomal RNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	8	20	Total	C	N	O	P	0	0
			431	192	80	139	20		

- Molecule 23 is a protein called Ribosomal protein L10a.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	B	213	Total	C	N	O		0	0
			1055	629	213	213			

- Molecule 24 is a protein called Ribosomal protein L36a.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	F	95	Total	C	N	O		0	0
			467	277	95	95			

- Molecule 25 is a RNA chain called tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	Y	75	Total	C	N	O	P	0	0
			1597	713	285	525	74		
25	V	75	Total	C	N	O	P	0	0
			1597	713	285	525	74		
25	W	76	Total	C	N	O	P	0	0
			1619	723	290	531	75		

- Molecule 26 is a RNA chain called mRNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	y	3	Total	C	N	O	P	0	0
			60	27	7	23	3		
26	v	3	Total	C	N	O	P	0	0
			60	27	7	23	3		

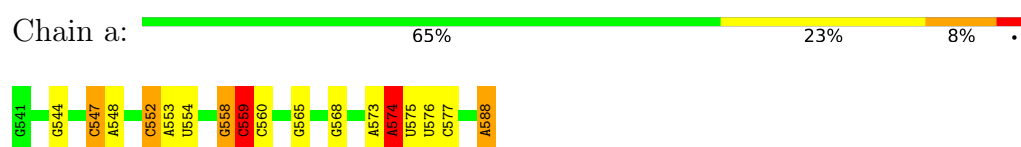
- Molecule 27 is a RNA chain called mRNA fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
27	w	2	44	20	10	12	2	0	0

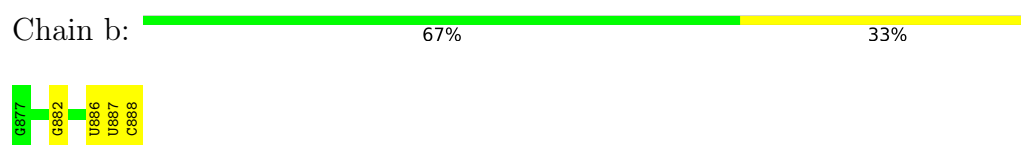
3 Residue-property plots [i](#)

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.

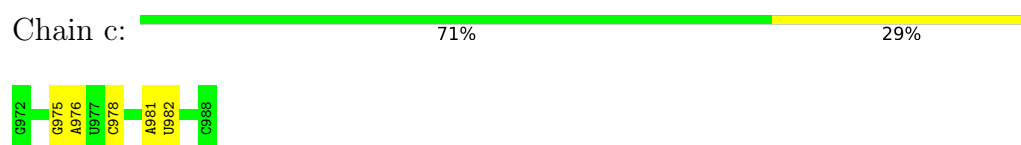
- Molecule 1: 40S ribosomal RNA fragment



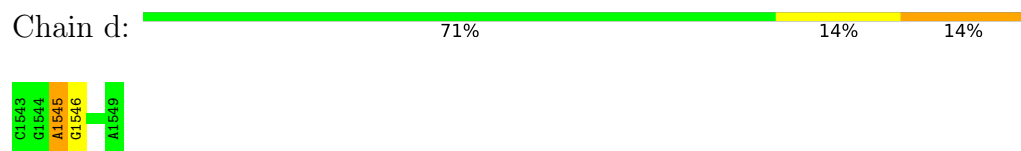
- Molecule 2: 40S ribosomal RNA fragment



- Molecule 3: 40S ribosomal RNA fragment



- Molecule 4: 40S ribosomal RNA fragment



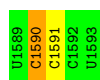
- Molecule 5: 40S ribosomal RNA fragment



There are no outlier residues recorded for this chain.

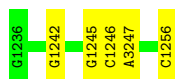
- Molecule 6: 40S ribosomal RNA fragment





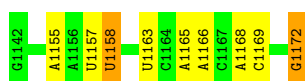
- Molecule 7: 40S ribosomal RNA fragment

Chain f: 76% 24%



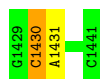
- Molecule 8: 40S ribosomal RNA fragment

Chain g: 71% 23% 6%



- Molecule 9: 40S ribosomal RNA fragment

Chain G: 85% 8% 8%



- Molecule 10: 40S ribosomal RNA fragment

Chain h: 88% 11%



- Molecule 11: Ribosomal protein S5

Chain T: 84% 15%




- Molecule 12: Ribosomal protein S14

Chain K: 79% 20%




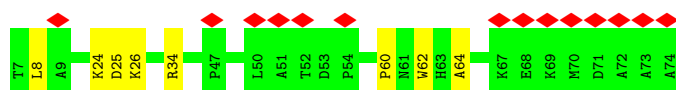
- Molecule 13: Ribosomal protein S23

Chain L:  81% 18%




- Molecule 14: Ribosomal protein S30

Chain X:  21% 88% 12%



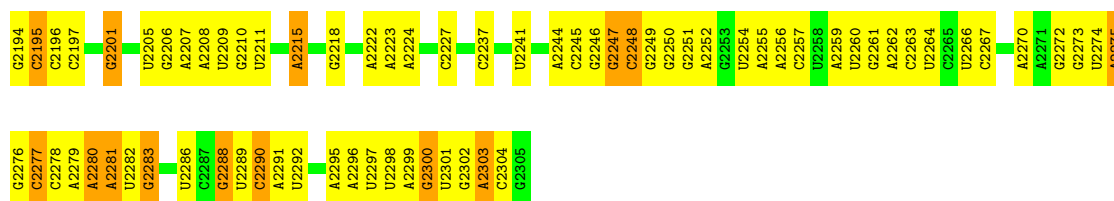
- Molecule 15: Ribosomal protein S15

Chain S:  91% 9%



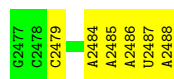
- Molecule 16: 60S ribosomal RNA fragment

Chain 2:  38% 50% 12%



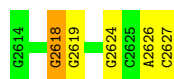
- Molecule 17: 60S ribosomal RNA fragment

Chain 3:  50% 50%



- Molecule 18: 60S ribosomal RNA fragment

Chain 4:  64% 29% 7%

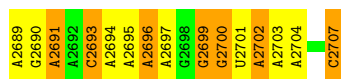
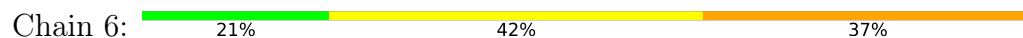


- Molecule 19: 60S ribosomal RNA fragment

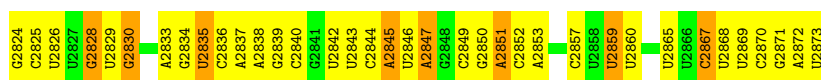
Chain 5:  17% 83%



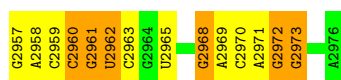
- Molecule 20: 60S ribosomal RNA fragment



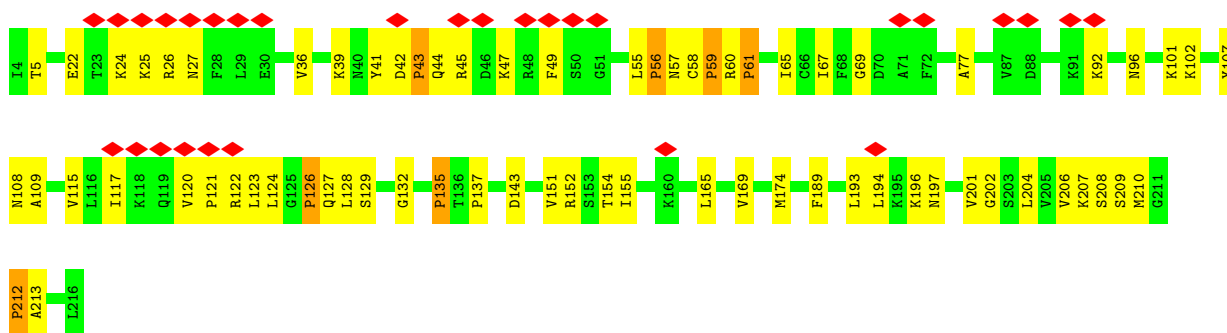
- Molecule 21: 60S ribosomal RNA fragment



- Molecule 22: 60S ribosomal RNA fragment



- Molecule 23: Ribosomal protein L10a



- Molecule 24: Ribosomal protein L36a



- Molecule 25: tRNA





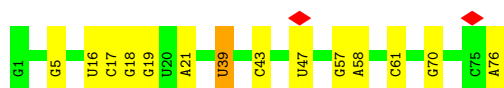
- Molecule 25: tRNA

Chain V: 79% 20% .



- Molecule 25: tRNA

Chain W: 82% 17% .



- Molecule 26: mRNA fragment

Chain y: 100%

There are no outlier residues recorded for this chain.

- Molecule 26: mRNA fragment

Chain v: 100%

There are no outlier residues recorded for this chain.

- Molecule 27: mRNA fragment

Chain w: 100%

There are no outlier residues recorded for this chain.

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	81946	Depositor
Resolution determination method	Not provided	
CTF correction method	CTF CORRECTION OF EACH DEFOCUS GROUP VOLUME PRIOR TO BACK PROJECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	4000	Depositor
Magnification	65520	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	21309.771	Depositor
Minimum map value	-5197.505	Depositor
Average map value	340.467	Depositor
Map value standard deviation	1535.379	Depositor
Recommended contour level	1000.0	Depositor
Map size (\AA)	453.6, 453.6, 453.6	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	2.52, 2.52, 2.52	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	a	0.75	1/1151 (0.1%)	1.00	7/1793 (0.4%)
2	b	0.55	0/291	0.79	0/452
3	c	0.66	0/404	0.92	1/627 (0.2%)
4	d	0.51	0/174	0.86	0/270
5	e	0.46	0/93	0.62	0/142
6	E	0.55	0/109	0.86	0/166
7	f	0.65	0/504	0.89	0/785
8	g	0.66	0/737	0.88	2/1146 (0.2%)
9	G	0.54	0/307	0.82	0/476
10	h	0.48	0/2650	0.75	1/4127 (0.0%)
11	T	0.45	0/1546	0.71	0/2079
12	K	0.42	0/1078	0.73	0/1452
13	L	0.41	0/1114	0.73	0/1485
14	X	0.36	0/566	0.70	0/753
15	S	0.38	0/1003	0.65	1/1342 (0.1%)
16	2	1.10	3/2677 (0.1%)	1.68	69/4170 (1.7%)
17	3	0.19	0/290	0.43	0/450
18	4	0.67	0/342	1.31	5/533 (0.9%)
19	5	0.69	0/141	1.32	1/217 (0.5%)
20	6	1.25	0/470	2.08	30/732 (4.1%)
21	7	1.06	2/1174 (0.2%)	2.34	33/1825 (1.8%)
22	8	1.45	4/482 (0.8%)	1.89	23/750 (3.1%)
23	B	0.34	0/1054	0.63	9/1468 (0.6%)
24	F	0.47	0/466	0.68	2/646 (0.3%)
25	V	0.44	0/1784	0.75	0/2780
25	W	0.43	0/1809	0.71	0/2819
25	Y	0.47	0/1784	0.74	0/2780
26	v	0.51	0/65	0.65	0/98
26	y	0.39	0/65	0.68	0/98
27	w	0.40	0/49	0.79	0/74
All	All	0.67	10/24379 (0.0%)	1.11	184/36535 (0.5%)

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
1	a	0	4
4	d	0	1
6	E	0	1
8	g	0	2
9	G	0	1
10	h	0	2
25	W	0	2
All	All	0	13

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
21	7	2845	A	C6-N1	-11.01	1.27	1.35
22	8	2969	A	N9-C4	-7.42	1.33	1.37
16	2	2283	G	N9-C8	6.14	1.42	1.37
22	8	2961	G	N9-C4	6.09	1.42	1.38
22	8	2958	A	N9-C4	-5.70	1.34	1.37
1	a	565	G	C2-N3	5.45	1.37	1.32
22	8	2968	G	C6-N1	-5.25	1.35	1.39
16	2	2302	G	C6-N1	-5.16	1.35	1.39
16	2	2279	A	N7-C5	-5.10	1.36	1.39
21	7	2830	G	N3-C4	-5.04	1.31	1.35

All (184) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
21	7	2845	A	N1-C6-N6	39.35	142.21	118.60
21	7	2845	A	C6-N1-C2	38.59	141.75	118.60
21	7	2845	A	C5-C6-N1	-34.02	100.69	117.70
21	7	2845	A	N1-C2-N3	-22.73	117.94	129.30
20	6	2689	A	C8-N9-C4	-13.86	100.26	105.80
21	7	2845	A	C5-C6-N6	-12.67	113.56	123.70
22	8	2958	A	C8-N9-C4	12.00	110.60	105.80
16	2	2195	C	N3-C4-C5	10.36	126.05	121.90
16	2	2289	U	C2-N3-C4	-10.03	120.98	127.00
20	6	2689	A	N7-C8-N9	9.96	118.78	113.80
16	2	2245	C	C6-N1-C2	-9.93	116.33	120.30
16	2	2283	G	C8-N9-C4	-9.78	102.49	106.40
16	2	2278	C	N1-C2-O2	9.43	124.56	118.90
22	8	2961	G	N3-C4-C5	-9.22	123.99	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	2	2277	C	C5-C6-N1	-8.95	116.53	121.00
10	h	1714	U	N1-C1'-C2'	8.92	125.60	114.00
21	7	2835	U	C2-N1-C1'	8.88	128.36	117.70
16	2	2247	G	C5-C6-O6	-8.79	123.33	128.60
20	6	2701	U	C5-C4-O4	-8.74	120.66	125.90
22	8	2961	G	N3-C4-N9	8.70	131.22	126.00
16	2	2302	G	N1-C6-O6	-8.67	114.70	119.90
21	7	2828	G	N1-C6-O6	8.48	124.99	119.90
21	7	2867	C	N1-C2-O2	-8.48	113.81	118.90
21	7	2828	G	C5-C6-O6	-8.42	123.55	128.60
20	6	2689	A	N9-C4-C5	8.36	109.14	105.80
16	2	2283	G	N3-C2-N2	-8.32	114.08	119.90
16	2	2247	G	C4-C5-N7	8.31	114.12	110.80
22	8	2958	A	N7-C8-N9	-8.27	109.67	113.80
21	7	2837	A	N1-C6-N6	-8.22	113.67	118.60
16	2	2303	A	C8-N9-C4	8.22	109.09	105.80
16	2	2290	C	N1-C2-O2	-7.81	114.22	118.90
16	2	2283	G	N1-C6-O6	7.75	124.55	119.90
21	7	2851	A	N1-C6-N6	-7.70	113.98	118.60
21	7	2835	U	C5-C4-O4	-7.64	121.32	125.90
16	2	2247	G	C6-C5-N7	-7.56	125.86	130.40
16	2	2277	C	C2-N3-C4	-7.54	116.13	119.90
16	2	2201	G	N1-C6-O6	7.51	124.40	119.90
18	4	2618	G	C5-C6-O6	7.50	133.10	128.60
16	2	2283	G	N3-C4-N9	-7.47	121.52	126.00
21	7	2835	U	N3-C4-O4	7.46	124.62	119.40
22	8	2957	G	N3-C4-N9	-7.43	121.54	126.00
16	2	2241	U	C5-C4-O4	7.39	130.33	125.90
1	a	588	A	N9-C1'-C2'	7.30	123.50	114.00
22	8	2960	C	C6-N1-C2	7.26	123.21	120.30
20	6	2689	A	N1-C2-N3	7.23	132.91	129.30
16	2	2302	G	C6-C5-N7	7.20	134.72	130.40
20	6	2700	G	C5-C6-O6	-7.13	124.32	128.60
16	2	2247	G	N1-C6-O6	7.10	124.16	119.90
16	2	2267	C	N1-C2-O2	-7.06	114.67	118.90
22	8	2957	G	N3-C4-C5	6.98	132.09	128.60
1	a	559	C	N1-C1'-C2'	6.92	123.00	114.00
3	c	981	A	N9-C1'-C2'	6.90	122.97	114.00
24	F	5	PRO	N-CA-CB	6.90	111.58	103.30
16	2	2248	C	N1-C2-O2	-6.86	114.78	118.90
22	8	2969	A	C2-N3-C4	-6.81	107.19	110.60
16	2	2277	C	N1-C2-O2	-6.79	114.83	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	2	2283	G	N7-C8-N9	6.77	116.48	113.10
21	7	2830	G	C5-C6-N1	-6.76	108.12	111.50
21	7	2824	G	C6-N1-C2	6.72	129.13	125.10
16	2	2304	C	C6-N1-C2	-6.68	117.63	120.30
20	6	2689	A	C4-N9-C1'	6.57	138.13	126.30
16	2	2278	C	N1-C2-N3	-6.56	114.61	119.20
16	2	2289	U	N3-C4-C5	6.52	118.51	114.60
21	7	2845	A	C2-N3-C4	-6.49	107.36	110.60
20	6	2700	G	C8-N9-C1'	-6.48	118.58	127.00
1	a	565	G	N3-C2-N2	6.46	124.43	119.90
16	2	2277	C	C6-N1-C2	6.46	122.89	120.30
20	6	2700	G	N3-C4-N9	6.41	129.85	126.00
23	B	56	PRO	N-CA-CB	6.39	110.97	103.30
8	g	1172	G	C2'-C3'-O3'	6.37	123.90	113.70
21	7	2847	A	C8-N9-C4	-6.35	103.26	105.80
21	7	2865	U	N1-C2-O2	-6.32	118.38	122.80
20	6	2689	A	N1-C6-N6	-6.31	114.81	118.60
16	2	2248	C	C6-N1-C2	6.30	122.82	120.30
16	2	2218	G	C5-C6-N1	6.30	114.65	111.50
21	7	2857	C	N3-C4-C5	6.25	124.40	121.90
20	6	2689	A	C4-C5-C6	6.25	120.12	117.00
16	2	2278	C	C6-N1-C1'	-6.18	113.38	120.80
16	2	2278	C	C2-N1-C1'	6.18	125.59	118.80
22	8	2968	G	N1-C6-O6	-6.16	116.20	119.90
21	7	2837	A	C8-N9-C4	6.14	108.26	105.80
20	6	2700	G	C4-N9-C1'	6.14	134.48	126.50
21	7	2837	A	N7-C8-N9	-6.12	110.74	113.80
16	2	2289	U	C5-C6-N1	-6.10	119.65	122.70
16	2	2280	A	C8-N9-C4	-6.09	103.36	105.80
23	B	61	PRO	N-CA-CB	6.05	110.56	103.30
22	8	2972	G	C8-N9-C1'	-6.05	119.13	127.00
16	2	2278	C	C2-N3-C4	6.05	122.92	119.90
1	a	574	A	N9-C1'-C2'	6.04	121.86	114.00
23	B	135	PRO	N-CA-CB	6.04	110.55	103.30
21	7	2869	U	N3-C2-O2	-6.00	118.00	122.20
8	g	1172	G	N9-C1'-C2'	5.99	121.79	114.00
16	2	2194	G	C8-N9-C4	5.99	108.80	106.40
23	B	121	PRO	N-CA-CB	5.99	110.49	103.30
20	6	2700	G	C6-C5-N7	-5.99	126.81	130.40
16	2	2283	G	C5-N7-C8	-5.98	101.31	104.30
20	6	2693	C	C6-N1-C1'	-5.98	113.62	120.80
21	7	2835	U	C6-N1-C1'	-5.98	112.83	121.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	2	2196	C	C6-N1-C2	-5.96	117.92	120.30
23	B	126	PRO	N-CA-CB	5.94	110.42	103.30
21	7	2868	U	N3-C2-O2	5.91	126.34	122.20
24	F	56	PRO	N-CA-CB	5.91	110.39	103.30
23	B	43	PRO	N-CA-CB	5.89	110.36	103.30
20	6	2701	U	N3-C4-O4	5.88	123.52	119.40
23	B	137	PRO	N-CA-CB	5.86	110.33	103.30
16	2	2303	A	N9-C4-C5	-5.85	103.46	105.80
20	6	2702	A	C8-N9-C4	-5.84	103.46	105.80
21	7	2867	C	N3-C2-O2	5.84	125.99	121.90
20	6	2693	C	C2-N1-C1'	5.84	125.22	118.80
23	B	59	PRO	N-CA-CB	5.84	110.30	103.30
22	8	2970	C	C2-N1-C1'	5.82	125.20	118.80
16	2	2201	G	C5-C6-O6	-5.81	125.11	128.60
23	B	212	PRO	N-CA-CB	5.81	110.27	103.30
16	2	2241	U	C6-N1-C1'	5.80	129.32	121.20
16	2	2302	G	C5-C6-N1	5.79	114.40	111.50
20	6	2696	A	C8-N9-C4	-5.79	103.48	105.80
20	6	2700	G	C4-C5-N7	5.78	113.11	110.80
16	2	2302	G	N3-C4-N9	-5.75	122.55	126.00
20	6	2689	A	N3-C4-C5	-5.74	122.78	126.80
20	6	2707	C	N3-C2-O2	5.72	125.90	121.90
21	7	2837	A	C4-C5-N7	-5.70	107.85	110.70
16	2	2283	G	N1-C2-N2	5.70	121.33	116.20
18	4	2618	G	C5-C6-N1	-5.69	108.65	111.50
16	2	2283	G	C2-N3-C4	-5.68	109.06	111.90
16	2	2290	C	C2-N3-C4	-5.67	117.06	119.90
16	2	2290	C	N3-C4-C5	5.66	124.16	121.90
16	2	2248	C	N3-C2-O2	5.64	125.85	121.90
22	8	2959	C	N1-C2-O2	-5.64	115.52	118.90
20	6	2707	C	C6-N1-C2	5.63	122.55	120.30
16	2	2196	C	N3-C4-C5	5.62	124.15	121.90
18	4	2618	G	N9-C4-C5	5.60	107.64	105.40
16	2	2301	U	C6-N1-C2	5.60	124.36	121.00
16	2	2197	C	N3-C2-O2	5.58	125.81	121.90
18	4	2618	G	C4-C5-N7	-5.56	108.57	110.80
1	a	558	G	N9-C1'-C2'	5.56	121.23	114.00
21	7	2853	A	N1-C6-N6	5.56	121.94	118.60
16	2	2300	G	C8-N9-C1'	-5.56	119.78	127.00
1	a	565	G	N1-C2-N2	-5.55	111.20	116.20
16	2	2301	U	N3-C2-O2	5.53	126.07	122.20
16	2	2241	U	N1-C2-N3	5.52	118.21	114.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
21	7	2830	G	C2-N3-C4	-5.52	109.14	111.90
16	2	2246	G	C5-C6-N1	5.51	114.26	111.50
16	2	2289	U	C5-C4-O4	-5.50	122.60	125.90
1	a	552	C	N1-C1'-C2'	5.47	121.11	114.00
16	2	2274	U	N3-C2-O2	-5.46	118.38	122.20
20	6	2707	C	N1-C2-O2	-5.44	115.63	118.90
16	2	2283	G	N3-C4-C5	5.42	131.31	128.60
16	2	2301	U	N1-C2-O2	-5.41	119.01	122.80
16	2	2195	C	C5-C4-N4	-5.40	116.42	120.20
22	8	2972	G	C4-N9-C1'	5.39	133.51	126.50
20	6	2700	G	N1-C6-O6	5.38	123.13	119.90
21	7	2869	U	C2-N1-C1'	5.38	124.15	117.70
22	8	2969	A	N3-C4-C5	5.36	130.56	126.80
20	6	2689	A	C5-C6-N6	5.35	127.98	123.70
20	6	2700	G	N9-C4-C5	-5.35	103.26	105.40
16	2	2194	G	N7-C8-N9	-5.34	110.43	113.10
16	2	2275	A	C5-C6-N1	-5.32	115.04	117.70
20	6	2702	A	N7-C8-N9	5.31	116.46	113.80
21	7	2859	U	N3-C2-O2	5.31	125.92	122.20
16	2	2247	G	C5-N7-C8	-5.31	101.65	104.30
22	8	2972	G	N1-C2-N3	5.31	127.08	123.90
16	2	2303	A	C5-C6-N1	5.30	120.35	117.70
20	6	2699	G	N3-C4-C5	5.29	131.24	128.60
15	S	131	THR	N-CA-C	5.27	125.24	111.00
22	8	2963	C	C2-N3-C4	-5.27	117.27	119.90
22	8	2960	C	C5-C6-N1	-5.26	118.37	121.00
16	2	2246	G	N3-C4-N9	5.24	129.15	126.00
22	8	2973	G	N7-C8-N9	-5.21	110.50	113.10
22	8	2968	G	C5-C6-N1	5.20	114.10	111.50
22	8	2962	U	C6-N1-C1'	5.17	128.44	121.20
20	6	2691	A	C2-N3-C4	5.15	113.18	110.60
16	2	2281	A	C4-C5-C6	5.13	119.57	117.00
22	8	2961	G	C4-N9-C1'	5.12	133.15	126.50
16	2	2215	A	C2-N3-C4	-5.10	108.05	110.60
18	4	2618	G	N3-C4-N9	-5.08	122.95	126.00
20	6	2707	C	C3'-C2'-C1'	5.08	105.56	101.50
21	7	2869	U	C6-N1-C2	-5.05	117.97	121.00
16	2	2201	G	C4-C5-N7	5.04	112.82	110.80
22	8	2973	G	C8-N9-C4	5.04	108.42	106.40
21	7	2828	G	C8-N9-C1'	-5.03	120.47	127.00
22	8	2970	C	C6-N1-C1'	-5.03	114.77	120.80
21	7	2851	A	C5-C6-N6	5.01	127.71	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	5	2653	C	C5-C4-N4	-5.01	116.69	120.20
16	2	2288	G	C6-N1-C2	-5.00	122.10	125.10

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
6	E	1590	C	Sidechain
9	G	1430	C	Sidechain
25	W	39	U	Sidechain
25	W	5	G	Sidechain
1	a	547	C	Sidechain
1	a	558	G	Sidechain
1	a	559	C	Sidechain
1	a	574	A	Sidechain
4	d	1545	A	Sidechain
8	g	1157	U	Sidechain
8	g	1158	U	Sidechain
10	h	1610	G	Sidechain
10	h	1714	U	Sidechain

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
11	T	190/192 (99%)	141 (74%)	33 (17%)	16 (8%)	1	12
12	K	138/140 (99%)	96 (70%)	28 (20%)	14 (10%)	0	9
13	L	139/141 (99%)	106 (76%)	17 (12%)	16 (12%)	0	6

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
14	X	66/68 (97%)	48 (73%)	12 (18%)	6 (9%)	1	11
15	S	123/125 (98%)	91 (74%)	25 (20%)	7 (6%)	1	18
23	B	211/213 (99%)	76 (36%)	67 (32%)	68 (32%)	0	0
24	F	93/95 (98%)	37 (40%)	28 (30%)	28 (30%)	0	0
All	All	960/974 (99%)	595 (62%)	210 (22%)	155 (16%)	0	3

All (155) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
11	T	56	PHE
11	T	60	GLN
11	T	63	ILE
11	T	75	GLY
11	T	158	PHE
12	K	70	SER
12	K	100	LYS
12	K	124	MET
12	K	142	ARG
13	L	45	ALA
13	L	62	GLN
13	L	87	PRO
13	L	88	MET
13	L	89	ASP
13	L	107	GLY
13	L	108	ARG
13	L	117	PRO
14	X	8	LEU
14	X	64	ALA
15	S	15	GLY
23	B	24	LYS
23	B	36	VAL
23	B	39	LYS
23	B	42	ASP
23	B	43	PRO
23	B	55	LEU
23	B	56	PRO
23	B	58	CYS
23	B	59	PRO
23	B	60	ARG
23	B	61	PRO
23	B	109	ALA

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Mol	Chain	Res	Type
23	B	115	VAL
23	B	120	VAL
23	B	122	ARG
23	B	124	LEU
23	B	126	PRO
23	B	127	GLN
23	B	128	LEU
23	B	129	SER
23	B	135	PRO
23	B	143	ASP
23	B	151	VAL
23	B	174	MET
23	B	201	VAL
23	B	207	LYS
23	B	208	SER
23	B	209	SER
23	B	210	MET
23	B	212	PRO
24	F	5	PRO
24	F	13	LYS
24	F	34	SER
24	F	47	GLN
24	F	53	GLN
24	F	66	LYS
24	F	93	LEU
11	T	58	LYS
11	T	100	GLY
11	T	157	ALA
11	T	159	LYS
11	T	183	SER
12	K	102	GLY
12	K	105	THR
12	K	110	PRO
12	K	139	SER
13	L	35	GLY
13	L	54	GLU
13	L	76	ARG
13	L	138	LYS
14	X	24	LYS
15	S	31	GLU
15	S	33	LEU
23	B	25	LYS

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Mol	Chain	Res	Type
23	B	26	ARG
23	B	41	TYR
23	B	44	GLN
23	B	47	LYS
23	B	67	ILE
23	B	77	ALA
23	B	96	ASN
23	B	117	ILE
23	B	123	LEU
23	B	169	VAL
23	B	193	LEU
23	B	194	LEU
23	B	196	LYS
23	B	197	ASN
23	B	204	LEU
23	B	206	VAL
24	F	6	LYS
24	F	9	LYS
24	F	11	TYR
24	F	21	THR
24	F	41	ARG
24	F	46	LYS
24	F	57	VAL
24	F	58	PHE
24	F	81	ALA
24	F	88	CYS
11	T	41	GLN
11	T	103	PRO
11	T	179	ASN
12	K	23	ALA
12	K	99	ALA
12	K	106	LYS
12	K	122	SER
12	K	141	ARG
13	L	109	GLN
14	X	60	PRO
15	S	130	PRO
23	B	22	GLU
23	B	49	PHE
23	B	57	ASN
23	B	101	LYS
23	B	102	LYS

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Mol	Chain	Res	Type
23	B	152	ARG
23	B	202	GLY
24	F	17	CYS
24	F	19	LYS
24	F	56	PRO
12	K	146	ARG
13	L	91	CYS
14	X	25	ASP
23	B	27	ASN
23	B	107	TYR
23	B	132	GLY
23	B	155	ILE
23	B	165	LEU
23	B	189	PHE
23	B	213	ALA
24	F	12	CYS
24	F	59	HIS
11	T	62	PRO
14	X	26	LYS
15	S	123	GLU
23	B	5	THR
23	B	45	ARG
23	B	92	LYS
23	B	108	ASN
23	B	154	THR
24	F	3	ASN
24	F	39	GLY
24	F	45	ARG
24	F	55	LYS
24	F	62	ALA
24	F	48	SER
11	T	102	ASN
11	T	138	PRO
15	S	104	GLY
23	B	69	GLY
13	L	110	GLY
23	B	65	ILE
13	L	114	GLY
15	S	53	GLY

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	
11	T	163/163 (100%)	147 (90%)	16 (10%)	8	26
12	K	112/112 (100%)	96 (86%)	16 (14%)	3	16
13	L	113/113 (100%)	101 (89%)	12 (11%)	6	24
14	X	57/57 (100%)	55 (96%)	2 (4%)	36	59
15	S	105/105 (100%)	102 (97%)	3 (3%)	42	64
All	All	550/550 (100%)	501 (91%)	49 (9%)	13	30

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
11	T	24	ILE
11	T	26	ASP
11	T	30	GLN
11	T	31	ASN
11	T	36	THR
11	T	41	GLN
11	T	44	VAL
11	T	51	TYR
11	T	56	PHE
11	T	57	ARG
11	T	66	ARG
11	T	73	PHE
11	T	74	HIS
11	T	117	ARG
11	T	139	MET
11	T	147	TYR
12	K	28	PHE
12	K	43	HIS
12	K	46	ASP
12	K	47	LEU
12	K	67	GLU
12	K	80	ASP
12	K	85	CYS

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Mol	Chain	Res	Type
12	K	95	ILE
12	K	113	GLN
12	K	116	LEU
12	K	117	ARG
12	K	119	LEU
12	K	129	ILE
12	K	140	THR
12	K	142	ARG
12	K	149	ARG
13	L	9	ILE
13	L	19	ARG
13	L	21	ASP
13	L	33	LEU
13	L	55	LYS
13	L	69	LYS
13	L	81	LYS
13	L	82	ILE
13	L	89	ASP
13	L	92	LEU
13	L	106	LEU
13	L	132	LEU
14	X	34	ARG
14	X	62	TRP
15	S	66	ARG
15	S	83	THR
15	S	124	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
11	T	10	GLN
11	T	46	HIS
11	T	52	GLN
11	T	96	HIS
11	T	102	ASN
11	T	143	ASN
11	T	199	ASN
12	K	43	HIS
13	L	78	ASN
13	L	98	ASN
14	X	42	ASN
15	S	29	ASN

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Mol	Chain	Res	Type
15	S	84	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	a	47/48 (97%)	15 (31%)	0
10	h	110/111 (99%)	12 (10%)	0
16	2	111/112 (99%)	54 (48%)	9 (8%)
17	3	11/12 (91%)	6 (54%)	1 (9%)
18	4	13/14 (92%)	5 (38%)	1 (7%)
19	5	5/6 (83%)	4 (80%)	1 (20%)
2	b	11/12 (91%)	4 (36%)	0
20	6	18/19 (94%)	13 (72%)	1 (5%)
21	7	49/50 (98%)	29 (59%)	5 (10%)
22	8	19/20 (95%)	8 (42%)	0
25	V	74/76 (97%)	15 (20%)	0
25	W	75/76 (98%)	13 (17%)	0
25	Y	74/76 (97%)	20 (27%)	1 (1%)
26	v	2/3 (66%)	0	0
26	y	2/3 (66%)	0	0
27	w	1/2 (50%)	0	0
3	c	16/17 (94%)	4 (25%)	0
4	d	6/7 (85%)	2 (33%)	0
5	e	3/4 (75%)	0	0
6	E	4/5 (80%)	1 (25%)	1 (25%)
7	f	20/21 (95%)	5 (25%)	0
8	g	30/31 (96%)	8 (26%)	0
9	G	12/13 (92%)	2 (16%)	1 (8%)
All	All	713/738 (96%)	220 (30%)	21 (2%)

All (220) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	a	544	G
1	a	547	C
1	a	548	A
1	a	552	C
1	a	553	A
1	a	554	U
1	a	559	C
1	a	560	C

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Mol	Chain	Res	Type
1	a	568	G
1	a	573	A
1	a	574	A
1	a	575	U
1	a	576	U
1	a	577	C
1	a	588	A
2	b	882	G
2	b	886	U
2	b	887	U
2	b	888	C
3	c	975	G
3	c	976	A
3	c	978	C
3	c	982	U
4	d	1545	A
4	d	1546	G
6	E	1591	C
7	f	1242	G
7	f	1245	G
7	f	1246	C
7	f	3247	A
7	f	1256	C
8	g	1155	A
8	g	1158	U
8	g	1163	U
8	g	1165	A
8	g	1166	A
8	g	1168	A
8	g	1169	C
8	g	1172	G
9	G	1430	C
9	G	1431	A
10	h	1608	C
10	h	1609	C
10	h	1624	G
10	h	1649	U
10	h	1651	G
10	h	1652	A
10	h	1661	G
10	h	1674	A
10	h	1710	G

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Mol	Chain	Res	Type
10	h	1713	G
10	h	1714	U
10	h	1715	A
16	2	2195	C
16	2	2201	G
16	2	2205	U
16	2	2206	G
16	2	2207	A
16	2	2208	A
16	2	2209	U
16	2	2210	G
16	2	2211	U
16	2	2215	A
16	2	2222	A
16	2	2223	A
16	2	2224	A
16	2	2227	C
16	2	2237	C
16	2	2244	A
16	2	2247	G
16	2	2248	C
16	2	2249	G
16	2	2250	G
16	2	2252	A
16	2	2254	U
16	2	2255	A
16	2	2256	A
16	2	2257	C
16	2	2259	A
16	2	2260	U
16	2	2261	G
16	2	2262	A
16	2	2263	C
16	2	2264	U
16	2	2266	U
16	2	2270	A
16	2	2272	G
16	2	2273	G
16	2	2275	A
16	2	2276	G
16	2	2277	C
16	2	2280	A

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Mol	Chain	Res	Type
16	2	2281	A
16	2	2282	U
16	2	2283	G
16	2	2286	U
16	2	2288	G
16	2	2290	C
16	2	2291	A
16	2	2292	U
16	2	2295	A
16	2	2296	A
16	2	2297	U
16	2	2298	U
16	2	2299	A
16	2	2300	G
16	2	2303	A
17	3	2479	C
17	3	2484	A
17	3	2485	A
17	3	2486	A
17	3	2487	U
17	3	2488	A
18	4	2618	G
18	4	2619	G
18	4	2624	G
18	4	2626	A
18	4	2627	C
19	5	2655	U
19	5	2656	A
19	5	2657	A
19	5	2658	G
20	6	2690	G
20	6	2691	A
20	6	2693	C
20	6	2694	A
20	6	2695	A
20	6	2696	A
20	6	2697	A
20	6	2699	G
20	6	2700	G
20	6	2702	A
20	6	2703	A
20	6	2704	A

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Mol	Chain	Res	Type
20	6	2707	C
21	7	2825	C
21	7	2826	U
21	7	2828	G
21	7	2829	U
21	7	2830	G
21	7	2833	A
21	7	2834	G
21	7	2835	U
21	7	2836	C
21	7	2838	A
21	7	2839	G
21	7	2840	C
21	7	2842	U
21	7	2843	U
21	7	2844	C
21	7	2845	A
21	7	2846	U
21	7	2847	A
21	7	2849	C
21	7	2850	G
21	7	2851	A
21	7	2852	C
21	7	2859	U
21	7	2860	U
21	7	2867	C
21	7	2870	C
21	7	2871	G
21	7	2872	A
21	7	2873	U
22	8	2960	C
22	8	2961	G
22	8	2962	U
22	8	2965	U
22	8	2968	G
22	8	2971	A
22	8	2972	G
22	8	2973	G
25	Y	2	C
25	Y	9	A
25	Y	16	U
25	Y	17	C

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Mol	Chain	Res	Type
25	Y	18	G
25	Y	19	G
25	Y	21	A
25	Y	23	A
25	Y	34	G
25	Y	42	C
25	Y	43	C
25	Y	47	U
25	Y	48	C
25	Y	52	G
25	Y	53	G
25	Y	70	G
25	Y	71	G
25	Y	72	C
25	Y	73	A
25	Y	74	C
25	V	8	U
25	V	17	C
25	V	18	G
25	V	19	G
25	V	20	U
25	V	21	A
25	V	22	G
25	V	42	C
25	V	46	G
25	V	47	U
25	V	48	C
25	V	52	G
25	V	61	C
25	V	69	G
25	V	75	C
25	W	16	U
25	W	17	C
25	W	18	G
25	W	19	G
25	W	21	A
25	W	39	U
25	W	43	C
25	W	47	U
25	W	57	G
25	W	58	A
25	W	61	C

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Mol	Chain	Res	Type
25	W	70	G
25	W	76	A

All (21) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
6	E	1590	C
9	G	1430	C
16	2	2211	U
16	2	2251	G
16	2	2254	U
16	2	2255	A
16	2	2270	A
16	2	2280	A
16	2	2281	A
16	2	2290	C
16	2	2297	U
17	3	2487	U
18	4	2618	G
19	5	2656	A
20	6	2699	G
21	7	2834	G
21	7	2850	G
21	7	2851	A
21	7	2859	U
21	7	2872	A
25	Y	15	G

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides 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.

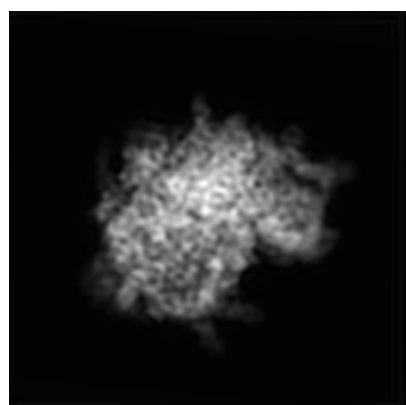
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-5327. These allow visual inspection of the internal detail of the map and identification of artifacts.

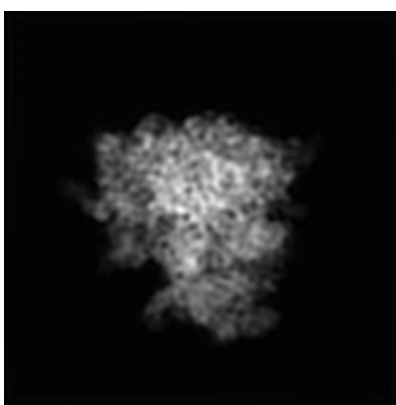
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

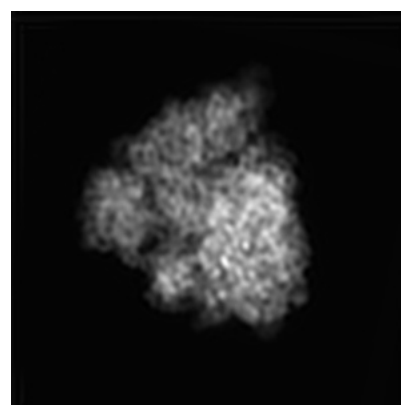
6.1.1 Primary map



X



Y

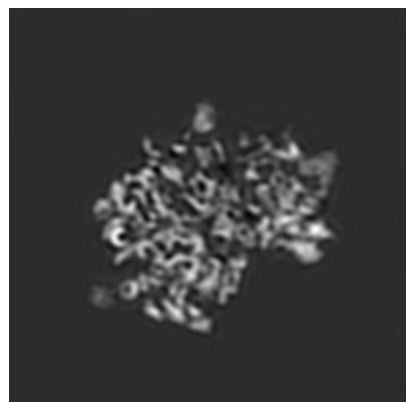


Z

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

6.2 Central slices [i](#)

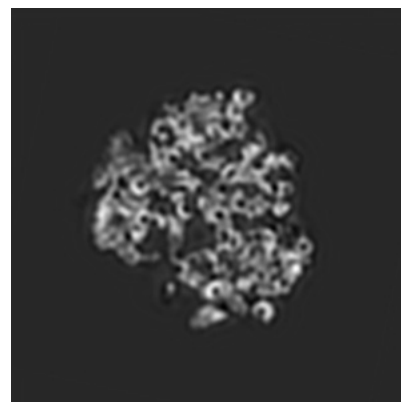
6.2.1 Primary map



X Index: 90



Y Index: 90

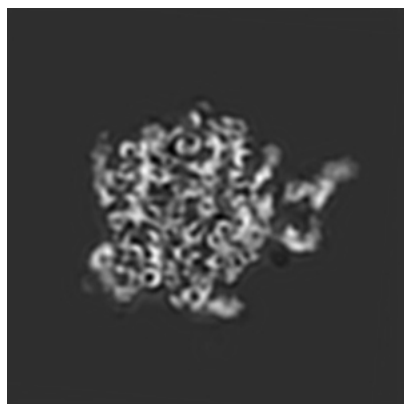


Z Index: 90

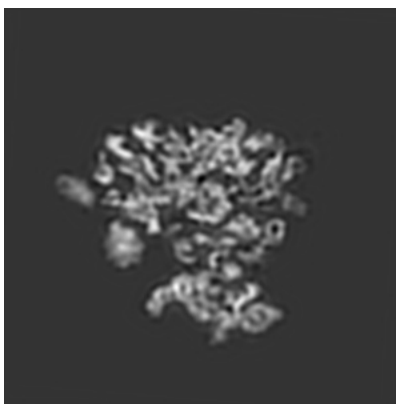
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

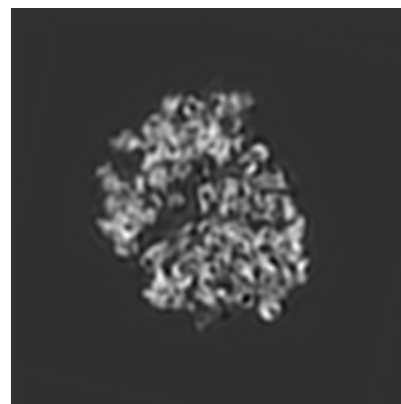
6.3.1 Primary map



X Index: 108



Y Index: 91



Z Index: 84

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

6.4 Orthogonal surface views [i](#)

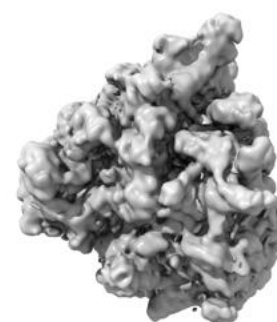
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 1000.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

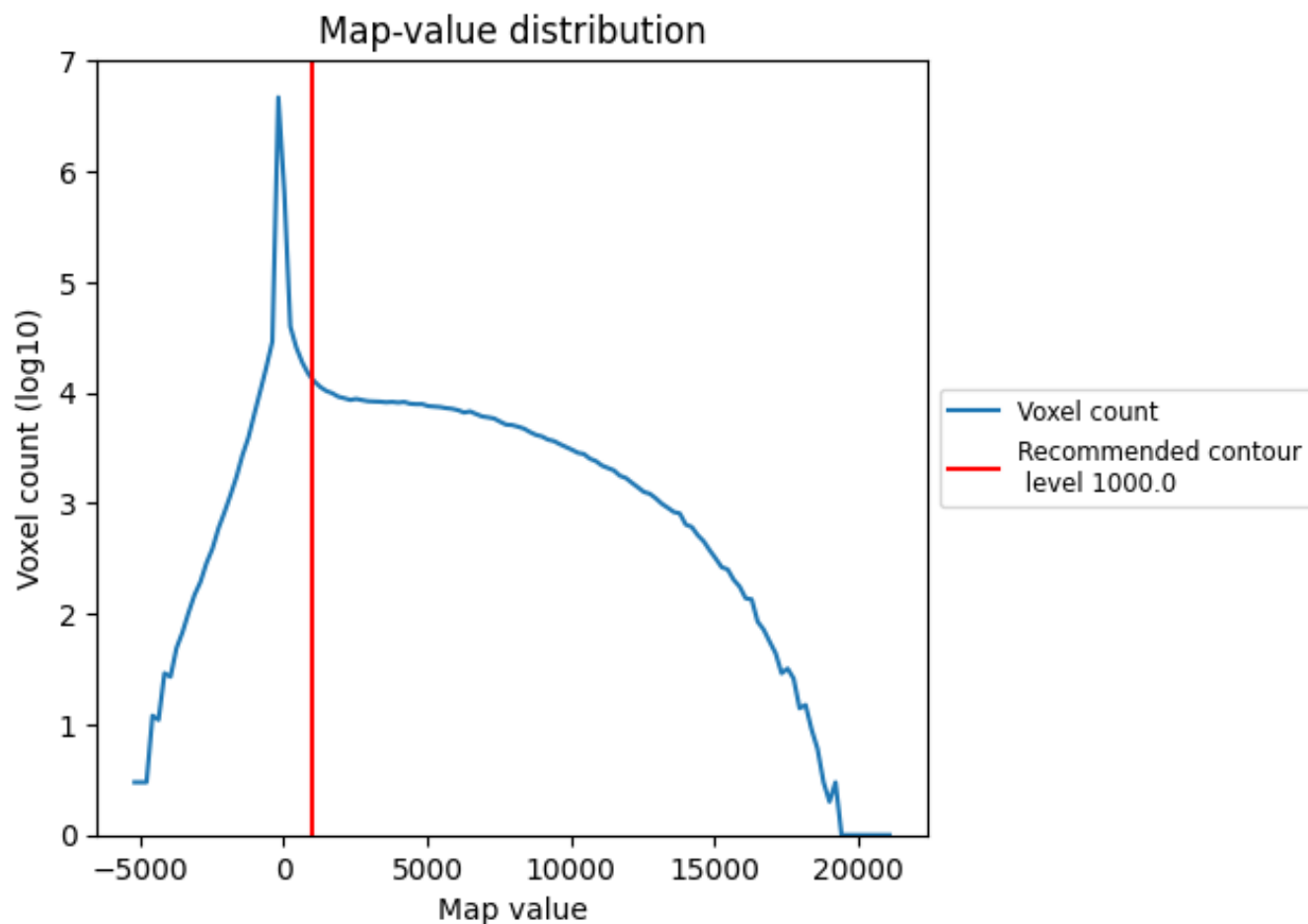
6.5 Mask visualisation

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

7 Map analysis [i](#)

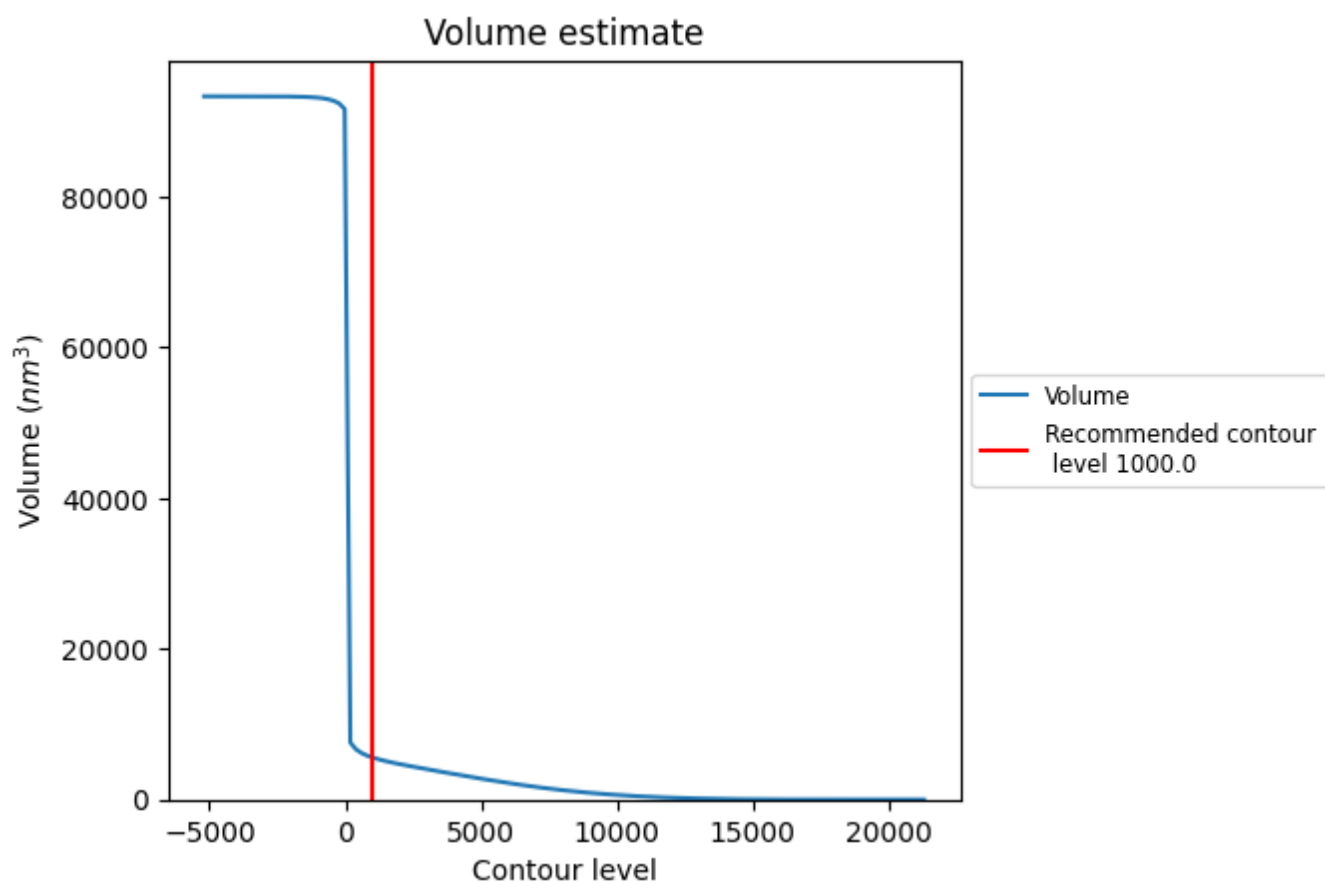
This section contains the results of statistical analysis of the map.

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.

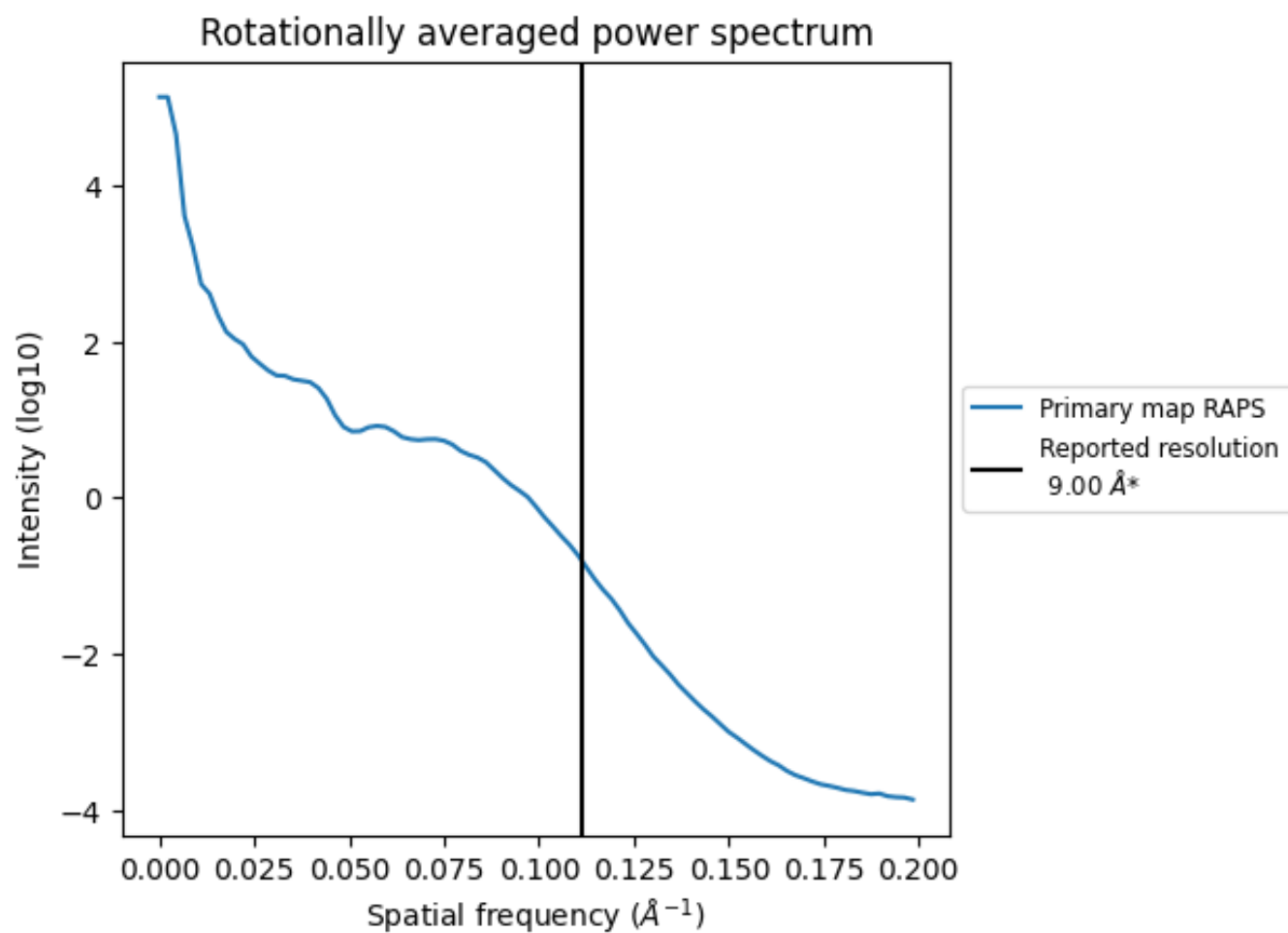
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 5581 nm^3 ; this corresponds to an approximate mass of 5042 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 0.111 Å⁻¹

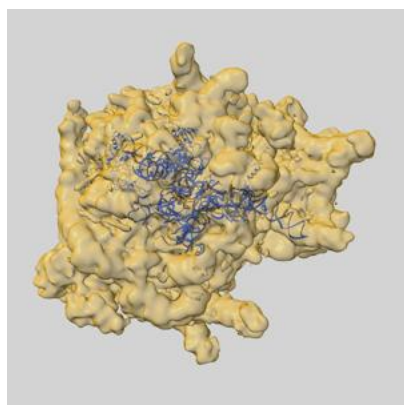
8 Fourier-Shell correlation ⓘ

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

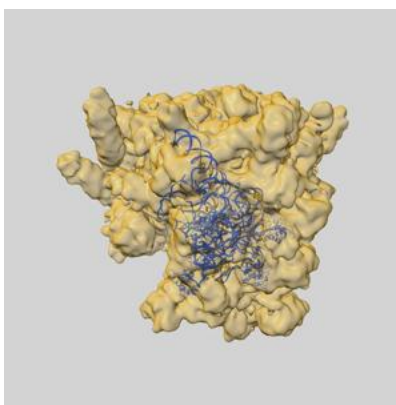
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-5327 and PDB model 3J0O. Per-residue inclusion information can be found in [section 3](#) on [page 9](#).

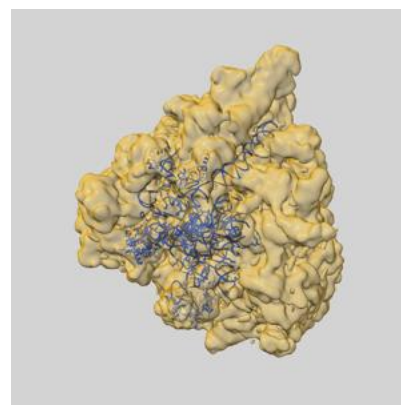
9.1 Map-model overlay [i](#)



X



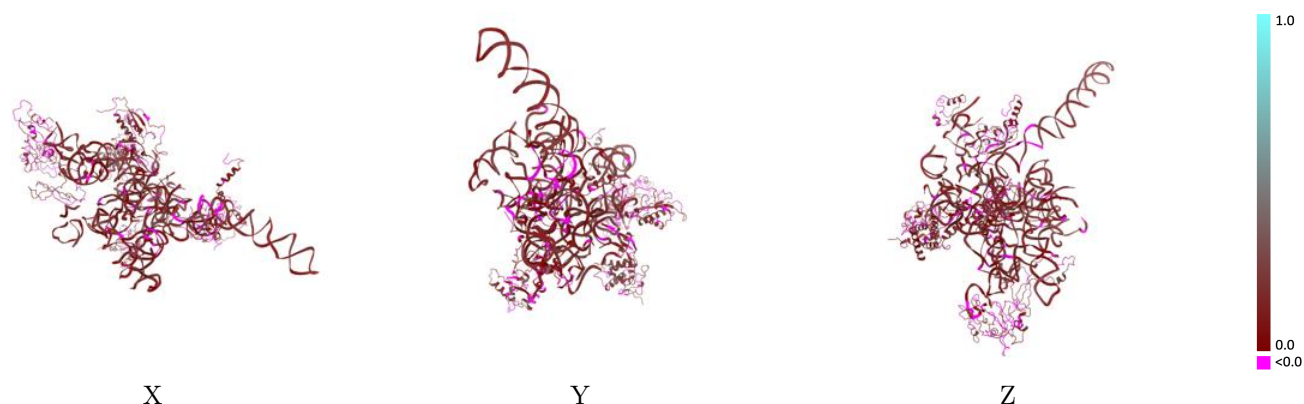
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 1000.0 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](#)



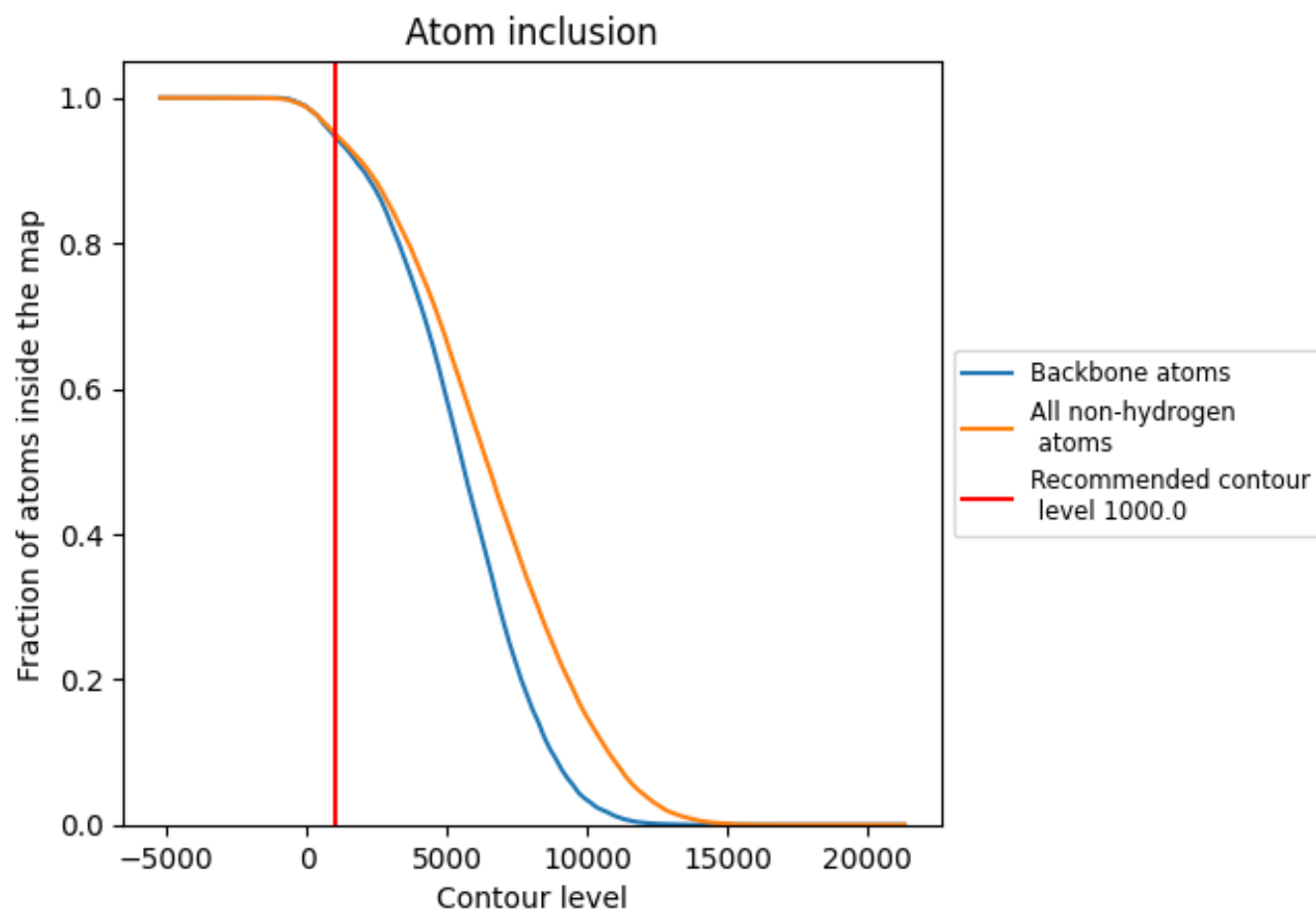
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1000.0).































































9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 95% 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 (1000.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9516	 0.1190
2	 0.9841	 0.1500
3	 1.0000	 0.1050
4	 0.9967	 0.1530
5	 1.0000	 0.1690
6	 1.0000	 0.1770
7	 0.9972	 0.1520
8	 0.9930	 0.1590
B	 0.8502	 0.0350
E	 1.0000	 0.1280
F	 0.9914	 0.1210
G	 1.0000	 0.1720
K	 0.9244	 0.0840
L	 0.9108	 0.0570
S	 0.9533	 0.0910
T	 0.9764	 0.1050
V	 0.9631	 0.1400
W	 0.9494	 0.1190
X	 0.7444	 0.0440
Y	 0.8773	 0.1200
a	 0.9961	 0.1350
b	 0.9962	 0.1440
c	 0.9862	 0.1360
d	 1.0000	 0.1320
e	 1.0000	 0.1310
f	 0.9978	 0.1380
g	 1.0000	 0.1320
h	 0.9451	 0.1310
v	 1.0000	 0.1860
w	 0.8636	 0.0260
y	 1.0000	 0.0860

