



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 12:58 AM UTC

PDB ID : 9FI8 / pdb_00009fi8
EMDB ID : EMD-50448
Title : SSU(head) structure derived from the SSU sample of the mitoribosome from *T. gondii*.
Authors : Rocha, R.E.O.; Barua, S.; Boissier, F.; Nguyen, T.T.; Hashem, Y.
Deposited on : 2024-05-28
Resolution : 3.60 Å(reported)
Based on initial model : .

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

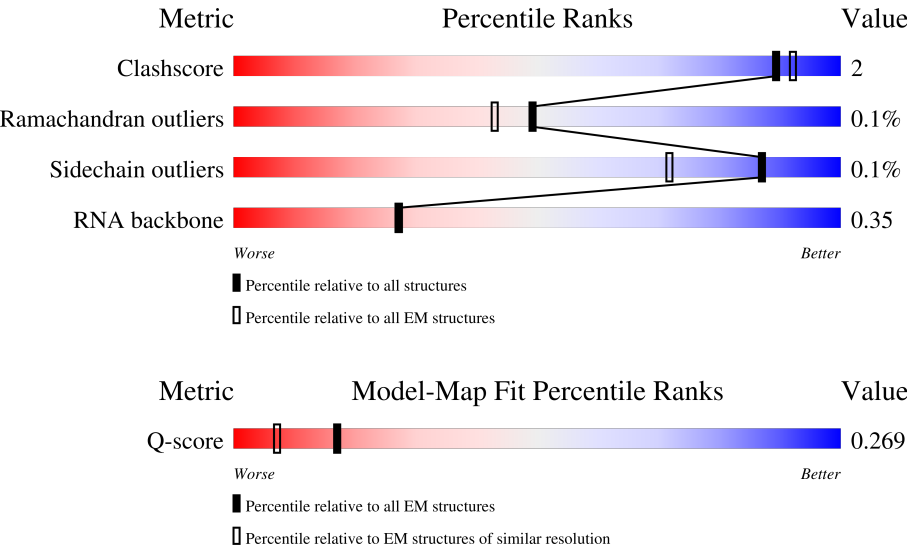
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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.



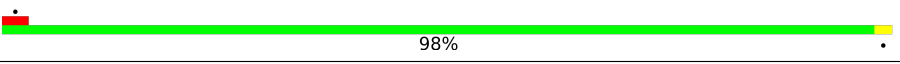
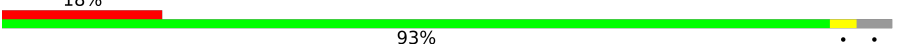
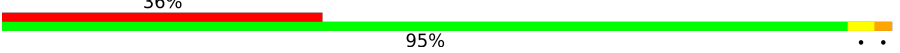
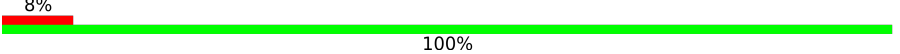
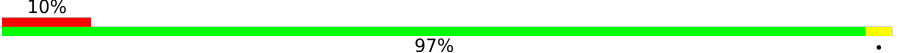
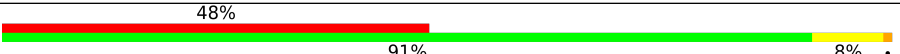
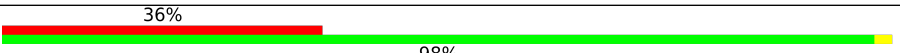
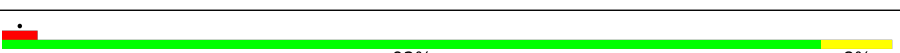
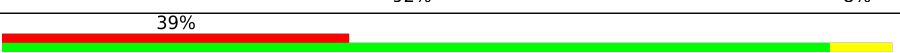
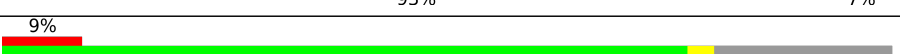
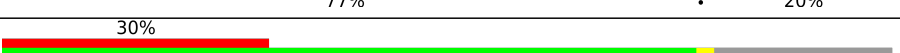
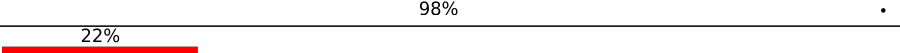
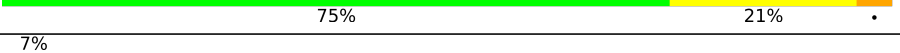
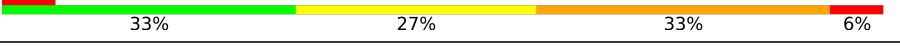
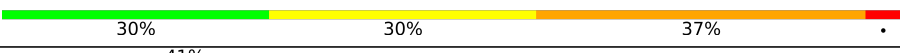
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	12797 (3.10 - 4.10)

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	HA	493	17% 79% 19%
2	HB	156	13% 95% 5%
3	HC	238	13% 74% 24%

Continued on next page...

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Mol	Chain	Length	Quality of chain
4	HD	102	
5	HE	177	
6	HF	354	
7	HG	97	
8	HH	90	
9	HI	144	
10	HJ	441	
11	HK	127	
12	HL	83	
13	HM	610	
14	HN	26	
15	HO	168	
16	HP	246	
17	HQ	757	
18	HR	170	
19	HS	121	
20	HT	405	
21	bN	58	
22	hB	56	
23	hD	15	
24	hG	68	
25	hL	33	
26	hM	10	
27	hR	27	
28	hS	17	

2 Entry composition

There are 28 unique types of molecules in this entry. The entry contains 40969 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 Small ribosomal subunit protein mS29,mS29(DAP3).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	HA	399	Total	C	N	O	S	0	0
			3287	2093	597	587	10		

- Molecule 2 is a protein called Ribosomal protein S9, putative.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	HB	156	Total	C	N	O	S	0	0
			1234	775	235	213	11		

- Molecule 3 is a protein called uS10m.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	HC	181	Total	C	N	O	S	0	0
			1432	905	259	262	6		

- Molecule 4 is a protein called Ribosomal protein S14, putative.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	HD	102	Total	C	N	O	S	0	0
			844	545	163	131	5		

- Molecule 5 is a protein called mS157.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	HE	119	Total	C	N	O	S	0	0
			926	581	184	154	7		

- Molecule 6 is a protein called DNA double-strand break repair rad50 ATPase.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	HF	341	Total	C	N	O	S	0	0
			2739	1730	513	489	7		

- Molecule 7 is a protein called mS149.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	HG	97	Total	C	N	O	S	0	0
			726	464	130	131	1		

- Molecule 8 is a protein called ms158.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	HH	90	Total	C	N	O	S	0	0
			761	500	122	136	3		

- Molecule 9 is a protein called mS151.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	HI	144	Total	C	N	O	S	0	0
			1203	787	216	194	6		

- Molecule 10 is a protein called 30S ribosomal protein S5, putative.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	HJ	430	Total	C	N	O	S	0	0
			3448	2211	585	645	7		

- Molecule 11 is a protein called mS155.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	HK	127	Total	C	N	O	S	0	0
			1045	660	198	182	5		

- Molecule 12 is a protein called mS159.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	HL	83	Total	C	N	O	S	0	0
			651	409	129	107	6		

- Molecule 13 is a protein called RAP domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	HM	525	Total	C	N	O	S	0	0
			4037	2552	728	744	13		

- Molecule 14 is a protein called mS161.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	HN	26	Total	C	N	O	S	0	0
			232	138	58	33	3		

- Molecule 15 is a protein called mS152.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	HO	168	Total	C	N	O	S	0	0
			1391	882	264	240	5		

- Molecule 16 is a protein called Small ribosomal subunit protein mS35 mitochondrial conserved domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	HP	197	Total	C	N	O	S	0	0
			1608	1018	301	284	5		

- Molecule 17 is a protein called mS141.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	HQ	607	Total	C	N	O	S	0	0
			4753	3014	843	880	16		

- Molecule 18 is a protein called AP2 domain transcription factor AP2IX-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	HR	162	Total	C	N	O	S	0	0
			1361	862	269	223	7		

- Molecule 19 is a protein called Acylphosphatase-like domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	HS	121	Total	C	N	O	S	0	0
			1001	629	186	181	5		

- Molecule 20 is a protein called mS33/mS27.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	HT	281	Total	C	N	O	S	0	0
			2313	1458	439	403	13		

- Molecule 21 is a RNA chain called SSUB-SSU.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	bN	58	Total	C	N	O	P	0	0
			1220	547	206	409	58		

- Molecule 22 is a RNA chain called SSUD.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	hB	56	Total	C	N	O	P	0	0
			1203	537	223	387	56		

- Molecule 23 is a RNA chain called ulr22.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	hD	15	Total	C	N	O	P	0	0
			300	135	30	120	15		

- Molecule 24 is a RNA chain called RNA9-SSU.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	hG	68	Total	C	N	O	P	0	0
			1436	643	246	479	68		

- Molecule 25 is a RNA chain called RNA5-SSU.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	hL	33	Total	C	N	O	P	0	0
			706	316	126	231	33		

- Molecule 26 is a RNA chain called ulr23.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	hM	10	Total	C	N	O	P	0	0
			200	90	20	80	10		

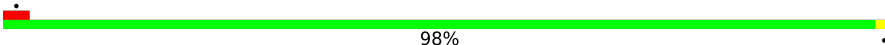
- Molecule 27 is a RNA chain called RNA19-SSU.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	hR	27	Total	C	N	O	P	0	0
			575	257	99	192	27		

- Molecule 28 is a RNA chain called ulr24.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	hS	17	Total	C	N	O	P	0	0
			337	153	34	134	16		

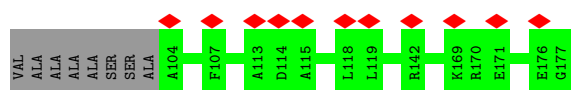
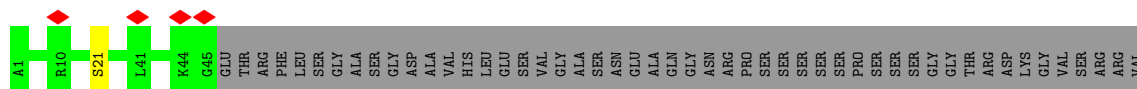
- Molecule 4: Ribosomal protein S14, putative

Chain HD:  98%

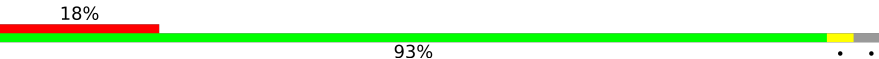


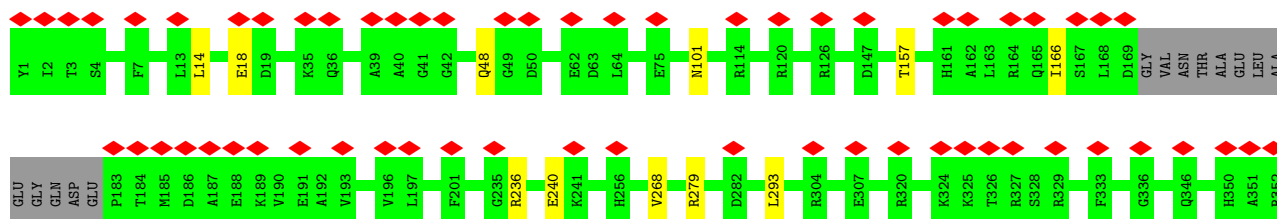
- Molecule 5: mS157

Chain HE:  67% 33%

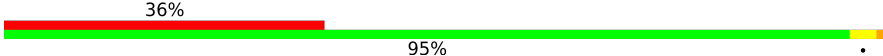


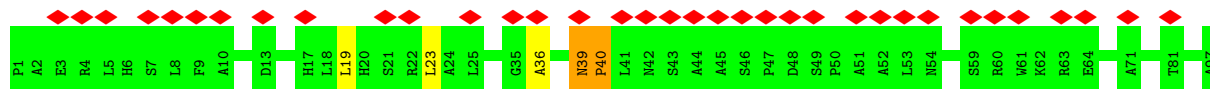
- Molecule 6: DNA double-strand break repair rad50 ATPase

Chain HF:  93%



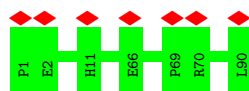
- Molecule 7: mS149

Chain HG:  95%

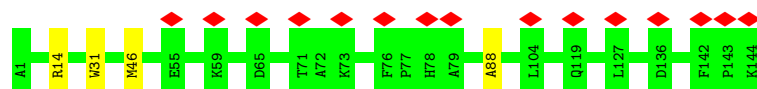


- Molecule 8: ms158

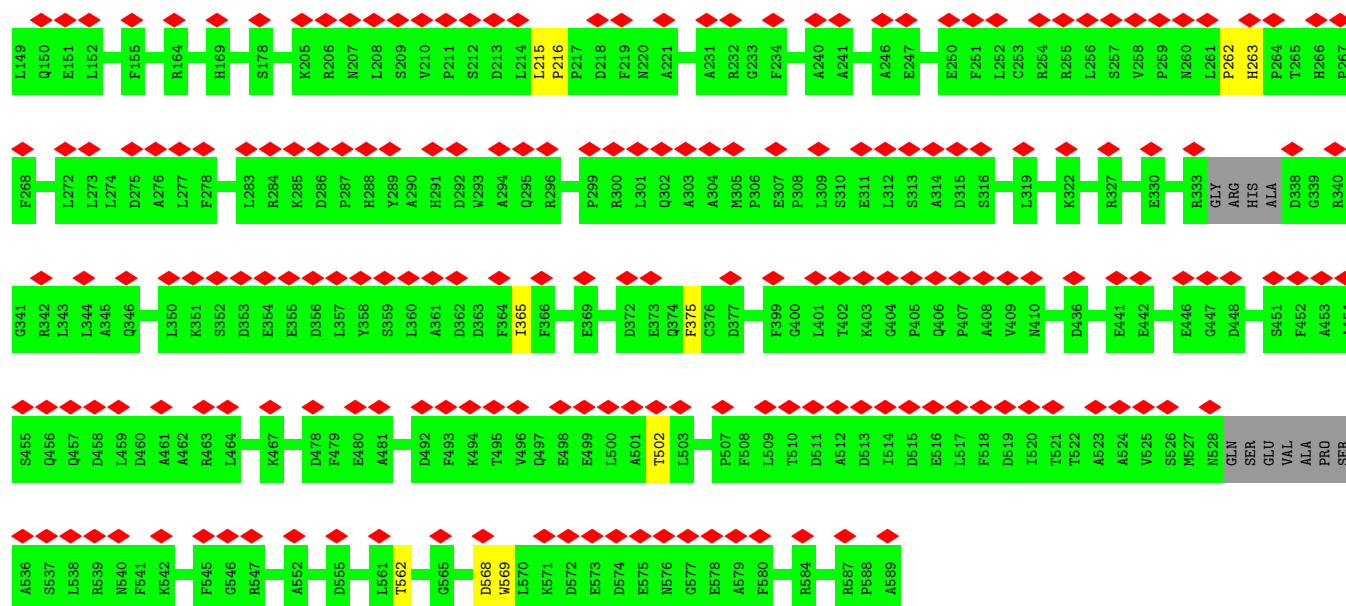
Chain HH:  100%



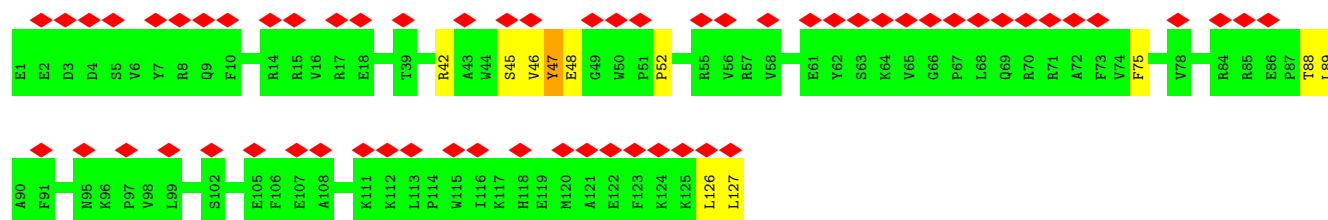
- Molecule 9: mS151



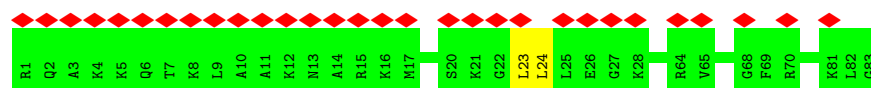
- Molecule 10: 30S ribosomal protein S5, putative



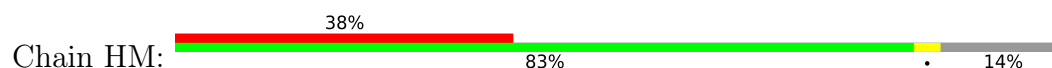
- Molecule 11: mS155

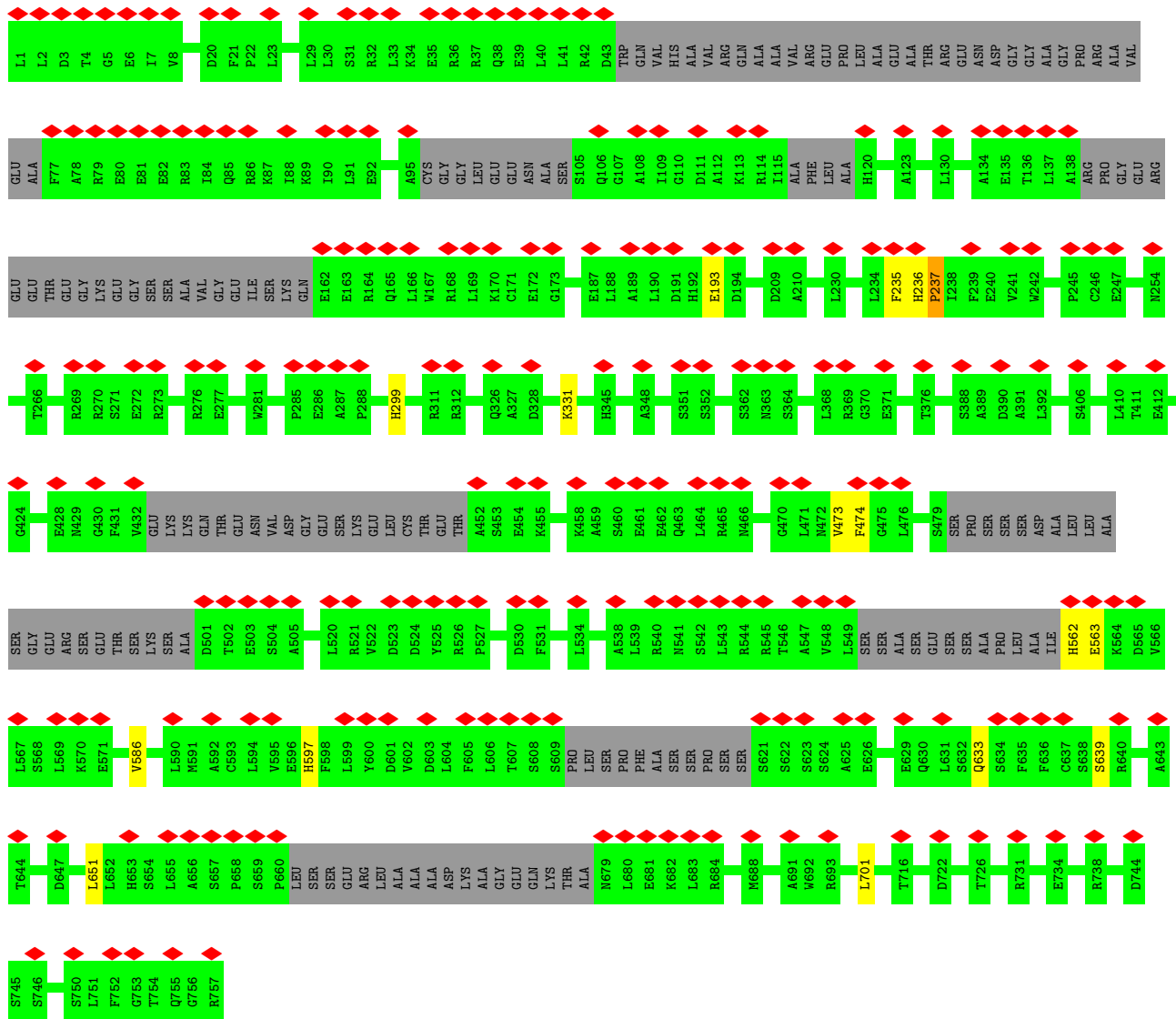


- Molecule 12: mS159

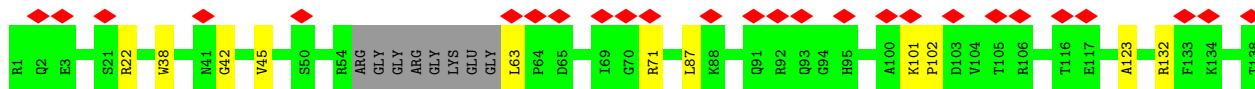
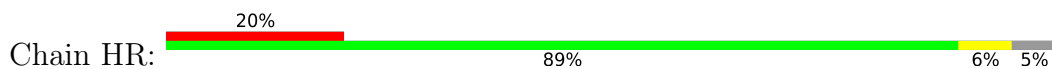


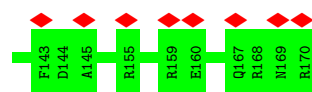
- Molecule 13: RAP domain-containing protein



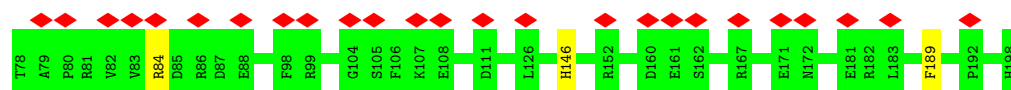


- Molecule 18: AP2 domain transcription factor AP2IX-6

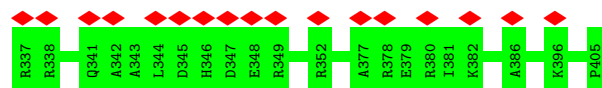
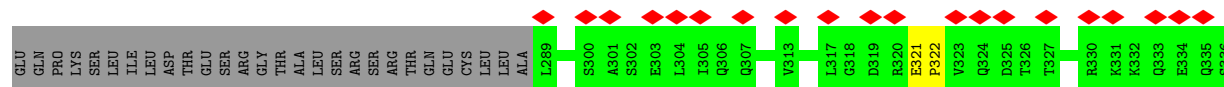
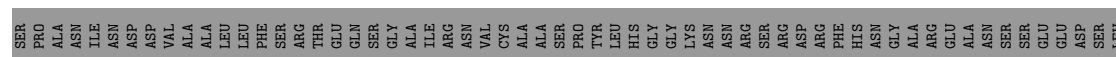
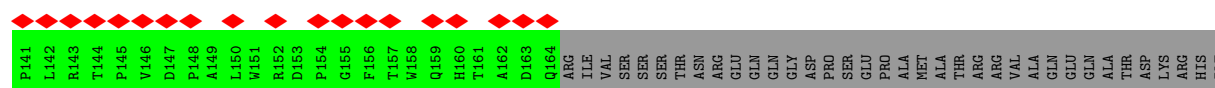
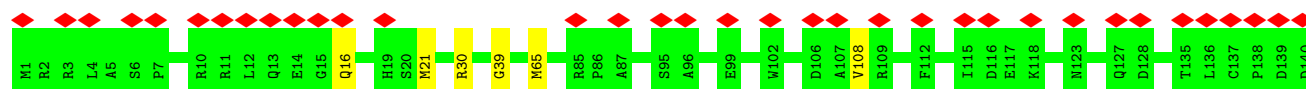




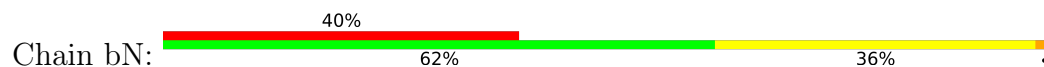
- Molecule 19: Acylphosphatase-like domain-containing protein



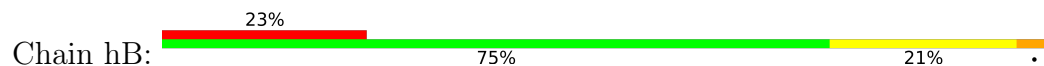
- Molecule 20: mS33/mS27



- Molecule 21: SSUB-SSU

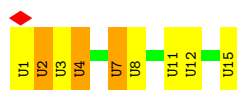


- Molecule 22: SSUD

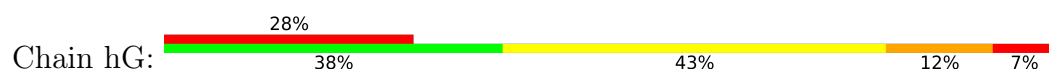


- Molecule 23: ulr22

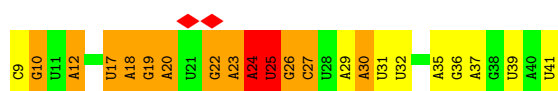




• Molecule 24: RNA9-SSU



• Molecule 25: RNA5-SSU



• Molecule 26: ulr23



• Molecule 27: RNA19-SSU



• Molecule 28: ulr24



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	22169	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	59000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	1.724	Depositor
Minimum map value	-0.016	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	632.394, 632.394, 632.394	wwPDB
Map dimensions	420, 420, 420	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.5057, 1.5057, 1.5057	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	HA	0.59	1/3359 (0.0%)	0.74	2/4529 (0.0%)
2	HB	0.62	0/1261	0.79	0/1699
3	HC	0.59	0/1473	0.73	0/1999
4	HD	0.71	0/867	0.91	3/1162 (0.3%)
5	HE	0.62	0/942	0.78	0/1259
6	HF	0.66	0/2795	0.82	1/3774 (0.0%)
7	HG	0.71	0/743	0.90	2/1016 (0.2%)
8	HH	0.57	0/789	0.74	0/1075
9	HI	0.65	0/1247	0.73	0/1683
10	HJ	0.59	0/3533	0.80	1/4794 (0.0%)
11	HK	0.71	0/1068	0.94	0/1433
12	HL	0.58	0/658	0.76	0/876
13	HM	0.67	0/4100	0.89	1/5556 (0.0%)
14	HN	0.80	0/232	1.10	0/299
15	HO	0.67	0/1423	0.96	5/1917 (0.3%)
16	HP	0.61	0/1644	0.79	2/2222 (0.1%)
17	HQ	0.68	1/4837 (0.0%)	0.87	2/6545 (0.0%)
18	HR	0.75	0/1395	0.93	0/1873
19	HS	0.65	0/1023	0.86	1/1371 (0.1%)
20	HT	0.64	0/2371	0.81	0/3198
21	bN	0.72	0/1360	0.79	0/2113
22	hB	0.71	0/1347	0.75	0/2099
23	hD	0.73	0/329	0.94	0/506
24	hG	0.92	2/1602 (0.1%)	1.29	22/2491 (0.9%)
25	hL	0.89	1/790 (0.1%)	1.34	13/1229 (1.1%)
26	hM	0.72	0/219	0.74	0/336
27	hR	0.79	0/642	1.08	4/998 (0.4%)
28	hS	0.97	0/370	1.58	8/570 (1.4%)
All	All	0.68	5/42419 (0.0%)	0.88	67/58622 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
24	hG	0	9
25	hL	0	18
26	hM	0	1
27	hR	0	15
28	hS	0	7
All	All	0	50

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
24	hG	40	A	O3'-P	12.83	1.80	1.61
24	hG	40	A	C3'-O3'	-8.76	1.29	1.42
25	hL	17	U	C3'-O3'	6.90	1.52	1.42
17	HQ	237	PRO	N-CD	-5.62	1.39	1.47
1	HA	423	ARG	CZ-NH1	-5.12	1.25	1.32

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	hG	40	A	C4'-C3'-O3'	15.58	136.37	113.00
25	hL	17	U	O3'-P-O5'	13.17	123.75	104.00
24	hG	56	G	C2'-C3'-O3'	12.21	127.81	109.50
24	hG	40	A	C2'-C3'-O3'	-11.54	96.39	113.70
24	hG	40	A	P-O3'-C3'	-11.46	103.01	120.20
25	hL	17	U	C4'-C3'-C2'	-11.24	91.36	102.60
24	hG	69	C	N1-C1'-C2'	10.39	127.58	112.00
25	hL	17	U	P-O3'-C3'	10.37	135.76	120.20
7	HG	39	ASN	OD1-CG-ND2	-8.68	113.92	122.60
17	HQ	237	PRO	N-CA-CB	8.63	110.85	103.25
15	HO	58	ARG	N-CA-C	8.22	119.86	111.07
24	hG	56	G	P-O3'-C3'	7.64	131.67	120.20
25	hL	31	U	O4'-C1'-N1	7.54	119.81	108.50
27	hR	6	U	N1-C1'-C2'	7.47	123.20	112.00
24	hG	57	G	P-O5'-C5'	-7.29	109.96	120.90
24	hG	40	A	C3'-C2'-O2'	7.21	121.52	110.70
24	hG	41	A	C4'-C3'-C2'	-7.20	95.41	102.60
10	HJ	568	ASP	CA-CB-CG	7.17	119.77	112.60
25	hL	25	U	O4'-C1'-N1	-7.12	97.53	108.20
25	hL	24	A	O4'-C1'-N9	7.04	118.77	108.20
25	hL	20	A	N9-C1'-C2'	6.91	122.37	112.00
28	hS	17	U	C2'-C3'-O3'	6.85	123.98	113.70
27	hR	8	U	N1-C1'-C2'	6.82	124.23	114.00
28	hS	15	U	O4'-C1'-N1	6.67	118.21	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	hG	41	A	O3'-P-O5'	6.59	113.88	104.00
24	hG	54	A	C4'-C3'-O3'	6.56	122.84	113.00
24	hG	40	A	C5'-C4'-C3'	-6.46	106.30	116.00
25	hL	17	U	O4'-C4'-C3'	6.41	110.41	104.00
24	hG	40	A	O3'-P-O5'	6.39	113.59	104.00
24	hG	48	U	C2'-C3'-O3'	6.34	119.02	109.50
28	hS	2	U	O4'-C1'-N1	6.04	117.57	108.50
28	hS	10	U	C4'-C3'-O3'	5.96	118.35	109.40
28	hS	7	U	O3'-P-O5'	5.90	112.85	104.00
27	hR	11	U	N1-C1'-C2'	5.87	120.80	112.00
19	HS	84	ARG	N-CA-C	5.80	119.94	112.86
25	hL	25	U	N1-C1'-C2'	5.79	122.69	114.00
28	hS	1	U	C1'-O4'-C4'	-5.76	104.14	109.90
28	hS	7	U	O4'-C1'-C2'	-5.74	100.06	105.80
15	HO	157	LYS	N-CA-C	5.71	118.12	110.35
24	hG	41	A	P-O5'-C5'	5.71	129.46	120.90
4	HD	4	PRO	N-CA-CB	5.69	106.38	103.19
13	HM	578	ILE	CB-CA-C	5.65	118.05	110.42
15	HO	154	VAL	N-CA-C	5.64	116.37	110.62
1	HA	423	ARG	NH1-CZ-NH2	-5.61	112.01	119.30
1	HA	423	ARG	NE-CZ-NH2	5.60	124.24	119.20
7	HG	40	PRO	N-CD-CG	5.59	111.58	103.20
25	hL	20	A	O4'-C1'-N9	-5.54	100.18	108.50
25	hL	30	A	C1'-C2'-O2'	-5.52	100.12	108.40
24	hG	66	G	C2'-C3'-O3'	-5.51	105.44	113.70
24	hG	65	U	C5'-C4'-C3'	-5.50	107.74	116.00
24	hG	44	C	O4'-C4'-C3'	5.46	109.46	104.00
24	hG	66	G	O5'-C5'-C4'	-5.45	103.33	111.50
15	HO	159	ASP	CA-C-N	5.40	126.59	119.84
15	HO	159	ASP	C-N-CA	5.40	126.59	119.84
24	hG	70	U	O4'-C1'-N1	5.40	116.30	108.20
16	HP	149	SER	CA-C-N	5.38	125.72	119.47
16	HP	149	SER	C-N-CA	5.38	125.72	119.47
25	hL	24	A	N9-C1'-C2'	-5.38	105.93	114.00
25	hL	17	U	C3'-C2'-C1'	5.37	106.67	101.30
17	HQ	597	HIS	CB-CG-CD2	-5.32	124.28	131.20
27	hR	18	A	O4'-C1'-N9	5.28	116.11	108.20
28	hS	8	U	C4'-C3'-C2'	-5.28	97.33	102.60
6	HF	48	GLN	OE1-CD-NE2	-5.27	117.33	122.60
24	hG	67	U	C3'-C2'-C1'	5.26	106.56	101.30
4	HD	4	PRO	O-C-N	5.21	123.71	121.31
24	hG	43	U	O4'-C1'-C2'	-5.10	102.50	107.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	HD	3	PRO	N-CA-CB	5.04	106.01	103.19

There are no chirality outliers.

All (50) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
24	hG	42	U	Sidechain
24	hG	43	U	Sidechain
24	hG	46	C	Sidechain
24	hG	47	U	Sidechain
24	hG	56	G	Sidechain
24	hG	63	U	Sidechain
24	hG	65	U	Sidechain
24	hG	66	G	Sidechain
24	hG	71	A	Sidechain
25	hL	10	G	Sidechain
25	hL	12	A	Sidechain
25	hL	17	U	Sidechain
25	hL	18	A	Sidechain
25	hL	19	G	Sidechain
25	hL	22	G	Sidechain
25	hL	23	A	Sidechain
25	hL	24	A	Sidechain
25	hL	25	U	Sidechain
25	hL	26	G	Sidechain
25	hL	27	C	Sidechain
25	hL	29	A	Sidechain
25	hL	32	U	Sidechain
25	hL	36	G	Sidechain
25	hL	37	A	Sidechain
25	hL	39	U	Sidechain
25	hL	41	U	Sidechain
25	hL	9	C	Sidechain
26	hM	3	U	Sidechain
27	hR	10	G	Sidechain
27	hR	11	U	Sidechain
27	hR	14	U	Sidechain
27	hR	15	A	Sidechain
27	hR	16	G	Sidechain
27	hR	17	C	Sidechain
27	hR	18	A	Sidechain
27	hR	19	U	Sidechain

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Mol	Chain	Res	Type	Group
27	hR	20	G	Sidechain
27	hR	21	G	Sidechain
27	hR	23	U	Sidechain
27	hR	24	U	Sidechain
27	hR	29	A	Sidechain
27	hR	32	A	Sidechain
27	hR	6	U	Sidechain
28	hS	1	U	Sidechain
28	hS	10	U	Sidechain
28	hS	11	U	Sidechain
28	hS	13	U	Sidechain
28	hS	15	U	Sidechain
28	hS	6	U	Sidechain
28	hS	7	U	Sidechain

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	HA	3287	0	3270	11	0
2	HB	1234	0	1234	9	0
3	HC	1432	0	1408	3	0
4	HD	844	0	864	0	0
5	HE	926	0	968	1	0
6	HF	2739	0	2791	8	0
7	HG	726	0	749	6	0
8	HH	761	0	733	0	0
9	HI	1203	0	1210	4	0
10	HJ	3448	0	3393	12	0
11	HK	1045	0	1062	7	0
12	HL	651	0	716	1	0
13	HM	4037	0	4171	12	0
14	HN	232	0	263	4	0
15	HO	1391	0	1414	8	0
16	HP	1608	0	1641	6	0
17	HQ	4753	0	4826	35	0
18	HR	1361	0	1356	6	0
19	HS	1001	0	970	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
20	HT	2313	0	2308	6	0
21	bN	1220	0	617	1	0
22	hB	1203	0	606	3	0
23	hD	300	0	151	15	0
24	hG	1436	0	729	10	0
25	hL	706	0	351	1	0
26	hM	200	0	101	9	0
27	hR	575	0	288	4	0
28	hS	337	0	172	1	0
All	All	40969	0	38362	143	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 2.

All (143) 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)
7:HG:39:ASN:OD1	7:HG:40:PRO:HD3	1.68	0.93
17:HQ:331:LYS:HZ2	23:hD:2:U:H5	0.95	0.93
17:HQ:193:GLU:HB3	17:HQ:235:PHE:CE2	2.10	0.86
2:HB:66:VAL:HG21	23:hD:7:U:N3	1.93	0.83
2:HB:66:VAL:HG21	23:hD:7:U:H3	1.43	0.83
11:HK:126:LEU:O	11:HK:127:LEU:OXT	1.96	0.83
2:HB:62:GLU:HB2	23:hD:7:U:O2	1.79	0.82
17:HQ:331:LYS:NZ	23:hD:2:U:H5	1.78	0.82
1:HA:423:ARG:NH1	26:hM:3:U:C5	2.47	0.81
11:HK:89:LEU:HD11	15:HO:125:ILE:HD11	1.64	0.78
17:HQ:235:PHE:CE2	17:HQ:237:PRO:HG2	2.18	0.78
1:HA:423:ARG:HH11	26:hM:3:U:H5	1.31	0.78
17:HQ:193:GLU:HB3	17:HQ:235:PHE:CZ	2.20	0.77
17:HQ:235:PHE:CE2	17:HQ:237:PRO:CG	2.69	0.76
13:HM:493:ASN:CB	13:HM:602:GLU:HG2	2.16	0.74
14:HN:8:LEU:HD21	14:HN:26:ARG:OXT	1.89	0.72
17:HQ:235:PHE:CE1	17:HQ:237:PRO:HD3	2.26	0.71
17:HQ:235:PHE:HE1	17:HQ:236:HIS:CE1	2.09	0.70
24:hG:40:A:O3'	24:hG:57:G:P	2.50	0.70
7:HG:36:ALA:HB3	10:HJ:569:TRP:CE3	2.29	0.68
7:HG:39:ASN:OD1	7:HG:40:PRO:CD	2.39	0.67
13:HM:493:ASN:HB2	13:HM:602:GLU:HG2	1.75	0.66
17:HQ:235:PHE:CG	17:HQ:237:PRO:HD2	2.30	0.66
17:HQ:235:PHE:CD2	17:HQ:237:PRO:HG2	2.33	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:HE:21:SER:HB2	26:hM:5:U:C5	2.34	0.63
6:HF:157:THR:HG21	15:HO:163:PHE:CE1	2.34	0.62
26:hM:6:U:H5'	26:hM:7:U:C5	2.34	0.62
13:HM:331:VAL:HB	13:HM:382:GLU:OE1	2.00	0.62
11:HK:75:PHE:CZ	11:HK:88:THR:HG21	2.34	0.61
10:HJ:365:ILE:CD1	10:HJ:375:PHE:CZ	2.84	0.60
17:HQ:193:GLU:HB3	17:HQ:235:PHE:CD2	2.37	0.60
17:HQ:235:PHE:CD1	17:HQ:237:PRO:HD3	2.37	0.59
17:HQ:299:HIS:CE1	23:hD:15:U:O4	2.54	0.59
10:HJ:365:ILE:HD13	10:HJ:375:PHE:CE2	2.37	0.59
17:HQ:193:GLU:CD	17:HQ:235:PHE:CE1	2.81	0.59
10:HJ:502:THR:O	10:HJ:502:THR:HG22	2.02	0.58
17:HQ:299:HIS:CE1	23:hD:15:U:C4	2.92	0.58
13:HM:493:ASN:HB3	13:HM:602:GLU:HG2	1.85	0.58
10:HJ:365:ILE:HD13	10:HJ:375:PHE:CZ	2.38	0.58
1:HA:244:ARG:CZ	26:hM:8:U:O2	2.53	0.57
17:HQ:235:PHE:CD2	17:HQ:237:PRO:CD	2.86	0.57
7:HG:36:ALA:HB3	10:HJ:569:TRP:CD2	2.39	0.57
3:HC:48:PRO:HD2	23:hD:15:U:O2	2.05	0.57
24:hG:8:A:N6	27:hR:30:A:C2	2.74	0.56
14:HN:8:LEU:CD2	14:HN:26:ARG:OXT	2.53	0.56
17:HQ:235:PHE:CZ	17:HQ:237:PRO:HD3	2.41	0.56
14:HN:8:LEU:CG	14:HN:26:ARG:OXT	2.54	0.55
9:HI:14:ARG:NH2	23:hD:11:U:C5	2.75	0.55
18:HR:22:ARG:CD	18:HR:87:LEU:HD13	2.36	0.55
17:HQ:235:PHE:CG	17:HQ:237:PRO:CD	2.89	0.55
24:hG:41:A:P	24:hG:57:G:OP2	2.64	0.55
17:HQ:586:VAL:HG21	17:HQ:651:LEU:HD21	1.88	0.55
15:HO:163:PHE:CE1	15:HO:167:LEU:HD23	2.42	0.54
17:HQ:193:GLU:HB3	17:HQ:235:PHE:CE1	2.43	0.54
9:HI:46:MET:SD	9:HI:88:ALA:O	2.66	0.54
17:HQ:235:PHE:CE1	17:HQ:236:HIS:CE1	2.94	0.54
13:HM:576:VAL:HG22	13:HM:577:SER:H	1.74	0.53
15:HO:163:PHE:CD1	15:HO:164:PRO:HA	2.45	0.52
9:HI:46:MET:HE1	9:HI:88:ALA:O	2.10	0.52
17:HQ:331:LYS:NZ	23:hD:2:U:C5	2.60	0.51
17:HQ:235:PHE:CD2	17:HQ:237:PRO:CG	2.95	0.50
11:HK:42:ARG:HD3	28:hS:6:U:C2	2.46	0.50
19:HS:189:PHE:HB2	27:hR:30:A:C2	2.46	0.49
10:HJ:365:ILE:HD11	10:HJ:375:PHE:CZ	2.46	0.49
14:HN:8:LEU:HD11	14:HN:26:ARG:OXT	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:HA:316:ARG:HD2	1:HA:328:GLU:OE1	2.11	0.49
15:HO:84:VAL:HG12	15:HO:84:VAL:O	2.12	0.49
1:HA:164:ILE:HG23	1:HA:174:VAL:HG11	1.96	0.48
2:HB:66:VAL:HG21	23:hD:7:U:C4	2.49	0.48
17:HQ:701:LEU:O	17:HQ:701:LEU:HG	2.13	0.48
17:HQ:235:PHE:CE2	17:HQ:237:PRO:CD	2.96	0.48
3:HC:8:ARG:NH2	3:HC:211:GLU:OE1	2.47	0.48
16:HP:59:ARG:NH2	25:hL:35:A:H61	2.12	0.47
17:HQ:235:PHE:CE1	17:HQ:236:HIS:ND1	2.82	0.47
17:HQ:235:PHE:CD2	17:HQ:237:PRO:HD2	2.48	0.47
1:HA:441:ARG:NH2	23:hD:4:U:C4	2.82	0.47
13:HM:220:VAL:HG12	13:HM:250:ILE:HD13	1.96	0.47
7:HG:36:ALA:CB	10:HJ:569:TRP:CE3	2.98	0.46
2:HB:62:GLU:CB	23:hD:7:U:O2	2.59	0.46
6:HF:14:LEU:HA	6:HF:18:GLU:OE1	2.15	0.46
15:HO:74:PHE:CZ	24:hG:43:U:H1'	2.51	0.46
16:HP:240:LEU:O	16:HP:244:VAL:HG23	2.16	0.45
24:hG:15:C:C5	27:hR:16:G:C6	3.04	0.45
6:HF:166:ILE:HG21	20:HT:108:VAL:HG12	1.97	0.45
17:HQ:235:PHE:CD1	17:HQ:237:PRO:CD	3.00	0.45
10:HJ:562:THR:O	10:HJ:569:TRP:CE3	2.70	0.45
6:HF:101:ASN:HD21	19:HS:146:HIS:HB3	1.82	0.44
24:hG:41:A:H5'	24:hG:57:G:OP2	2.17	0.44
6:HF:236:ARG:HG3	6:HF:240:GLU:CD	2.43	0.44
11:HK:75:PHE:CZ	11:HK:88:THR:CG2	3.01	0.44
17:HQ:473:VAL:HG12	17:HQ:474:PHE:H	1.83	0.44
22:hB:27:U:OP2	22:hB:52:G:N2	2.51	0.44
24:hG:41:A:P	24:hG:57:G:P	3.15	0.44
1:HA:244:ARG:NE	26:hM:8:U:O2	2.51	0.44
16:HP:202:ASP:O	20:HT:30:ARG:NH1	2.51	0.44
26:hM:6:U:H4'	26:hM:7:U:H6	1.83	0.44
3:HC:107:VAL:HG13	3:HC:113:VAL:HG21	1.99	0.43
9:HI:31:TRP:CD1	16:HP:172:VAL:HG21	2.52	0.43
18:HR:22:ARG:HD3	18:HR:87:LEU:HD13	1.99	0.43
24:hG:40:A:O2'	24:hG:41:A:P	2.75	0.43
18:HR:38:TRP:CE2	18:HR:45:VAL:HG23	2.53	0.43
10:HJ:215:LEU:HG	10:HJ:216:PRO:HD3	2.00	0.43
13:HM:320:THR:HG21	13:HM:360:LEU:HB2	2.00	0.43
17:HQ:193:GLU:O	17:HQ:235:PHE:CE2	2.72	0.43
13:HM:565:CYS:SG	13:HM:566:LYS:N	2.92	0.42
15:HO:111:SER:O	15:HO:115:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:HK:126:LEU:C	11:HK:127:LEU:OXT	2.62	0.42
17:HQ:299:HIS:NE2	23:hD:15:U:O4	2.52	0.42
13:HM:121:HIS:HB3	13:HM:122:PRO:HD3	2.00	0.42
20:HT:16:GLN:NE2	20:HT:21:MET:SD	2.92	0.42
12:HL:23:LEU:HD23	12:HL:24:LEU:N	2.34	0.42
18:HR:123:ALA:O	18:HR:132:ARG:NH2	2.52	0.42
16:HP:70:ARG:NH1	20:HT:39:GLY:O	2.52	0.42
2:HB:109:ARG:NH1	2:HB:130:ASP:OD2	2.52	0.42
17:HQ:562:HIS:O	17:HQ:563:GLU:HG3	2.19	0.42
22:hB:43:C:H1'	24:hG:66:G:H1'	2.02	0.42
6:HF:279:ARG:NH2	16:HP:221:ASP:OD1	2.53	0.42
11:HK:47:TYR:CG	11:HK:48:GLU:N	2.88	0.41
17:HQ:193:GLU:OE2	17:HQ:235:PHE:CD1	2.72	0.41
1:HA:244:ARG:HD3	26:hM:8:U:N3	2.36	0.41
2:HB:151:ARG:NH2	22:hB:24:U:OP1	2.53	0.41
17:HQ:235:PHE:CD1	17:HQ:236:HIS:N	2.88	0.41
6:HF:157:THR:HG21	15:HO:163:PHE:CD1	2.54	0.41
7:HG:19:LEU:CD1	7:HG:23:LEU:HB2	2.51	0.41
18:HR:101:LYS:N	18:HR:102:PRO:HD2	2.34	0.41
23:hD:1:U:O2'	23:hD:2:U:OP2	2.31	0.41
1:HA:316:ARG:HB2	1:HA:328:GLU:OE1	2.20	0.41
1:HA:426:ARG:NH1	26:hM:3:U:O2	2.54	0.41
10:HJ:215:LEU:N	10:HJ:216:PRO:CD	2.83	0.41
17:HQ:633:GLN:NE2	17:HQ:639:SER:OG	2.54	0.41
21:bN:101:A:O2'	21:bN:105:U:OP2	2.35	0.41
24:hG:40:A:H2'	24:hG:41:A:OP2	2.21	0.41
1:HA:298:LEU:HD11	13:HM:168:GLN:OE1	2.21	0.41
2:HB:67:ARG:NH1	2:HB:104:GLN:OE1	2.54	0.41
2:HB:136:ASP:OD1	2:HB:136:ASP:N	2.54	0.41
27:hR:29:A:H2'	27:hR:30:A:C8	2.55	0.41
18:HR:63:LEU:HD12	18:HR:71:ARG:HH22	1.85	0.40
13:HM:252:GLU:HB2	13:HM:255:ARG:HA	2.03	0.40
20:HT:65:MET:O	20:HT:65:MET:HG2	2.21	0.40
13:HM:282:LEU:HD12	13:HM:283:LEU:HD12	2.02	0.40
10:HJ:262:PRO:O	10:HJ:263:HIS:CD2	2.74	0.40
6:HF:268:VAL:HG23	6:HF:293:LEU:HD23	2.04	0.40
20:HT:321:GLU:N	20:HT:322:PRO:CD	2.85	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	HA	391/493 (79%)	377 (96%)	14 (4%)	0	100	100
2	HB	154/156 (99%)	145 (94%)	9 (6%)	0	100	100
3	HC	175/238 (74%)	159 (91%)	16 (9%)	0	100	100
4	HD	100/102 (98%)	95 (95%)	5 (5%)	0	100	100
5	HE	115/177 (65%)	108 (94%)	7 (6%)	0	100	100
6	HF	337/354 (95%)	315 (94%)	22 (6%)	0	100	100
7	HG	95/97 (98%)	89 (94%)	6 (6%)	0	100	100
8	HH	88/90 (98%)	87 (99%)	1 (1%)	0	100	100
9	HI	142/144 (99%)	136 (96%)	6 (4%)	0	100	100
10	HJ	424/441 (96%)	390 (92%)	34 (8%)	0	100	100
11	HK	125/127 (98%)	109 (87%)	14 (11%)	2 (2%)	7	36
12	HL	81/83 (98%)	76 (94%)	5 (6%)	0	100	100
13	HM	517/610 (85%)	465 (90%)	52 (10%)	0	100	100
14	HN	24/26 (92%)	21 (88%)	3 (12%)	0	100	100
15	HO	166/168 (99%)	153 (92%)	13 (8%)	0	100	100
16	HP	193/246 (78%)	183 (95%)	10 (5%)	0	100	100
17	HQ	587/757 (78%)	547 (93%)	40 (7%)	0	100	100
18	HR	158/170 (93%)	152 (96%)	5 (3%)	1 (1%)	21	54
19	HS	119/121 (98%)	112 (94%)	7 (6%)	0	100	100
20	HT	277/405 (68%)	259 (94%)	18 (6%)	0	100	100
All	All	4268/5005 (85%)	3978 (93%)	287 (7%)	3 (0%)	49	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
11	HK	46	VAL

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Mol	Chain	Res	Type
11	HK	52	PRO
18	HR	42	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	
1	HA	346/424 (82%)	345 (100%)	1 (0%)	86	83
2	HB	125/125 (100%)	125 (100%)	0	100	100
3	HC	157/206 (76%)	157 (100%)	0	100	100
4	HD	84/85 (99%)	84 (100%)	0	100	100
5	HE	97/140 (69%)	97 (100%)	0	100	100
6	HF	293/303 (97%)	293 (100%)	0	100	100
7	HG	77/77 (100%)	77 (100%)	0	100	100
8	HH	81/81 (100%)	81 (100%)	0	100	100
9	HI	127/127 (100%)	127 (100%)	0	100	100
10	HJ	372/380 (98%)	372 (100%)	0	100	100
11	HK	110/110 (100%)	108 (98%)	2 (2%)	51	68
12	HL	68/68 (100%)	68 (100%)	0	100	100
13	HM	431/496 (87%)	431 (100%)	0	100	100
14	HN	24/24 (100%)	24 (100%)	0	100	100
15	HO	149/149 (100%)	149 (100%)	0	100	100
16	HP	175/214 (82%)	175 (100%)	0	100	100
17	HQ	520/634 (82%)	520 (100%)	0	100	100
18	HR	137/141 (97%)	137 (100%)	0	100	100
19	HS	107/107 (100%)	107 (100%)	0	100	100
20	HT	242/346 (70%)	242 (100%)	0	100	100
All	All	3722/4237 (88%)	3719 (100%)	3 (0%)	87	87

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	HA	348	LEU
11	HK	45	SER
11	HK	47	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	HA	161	ASN
1	HA	302	HIS
1	HA	351	GLN
2	HB	79	GLN
6	HF	6	ASN
6	HF	36	GLN
7	HG	42	ASN
9	HI	16	HIS
9	HI	37	HIS
9	HI	108	ASN
10	HJ	263	HIS
10	HJ	288	HIS
10	HJ	346	GLN
10	HJ	506	ASN
10	HJ	576	ASN
11	HK	76	ASN
12	HL	49	GLN
12	HL	57	HIS
13	HM	264	ASN
13	HM	272	GLN
13	HM	353	GLN
13	HM	373	GLN
15	HO	25	ASN
16	HP	77	HIS
17	HQ	125	ASN
17	HQ	342	GLN
18	HR	19	GLN
19	HS	109	ASN
19	HS	172	ASN
20	HT	307	GLN
20	HT	358	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
21	bN	57/58 (98%)	21 (36%)	0
22	hB	55/56 (98%)	12 (21%)	0
23	hD	14/15 (93%)	6 (42%)	0
24	hG	67/68 (98%)	31 (46%)	0
25	hL	32/33 (96%)	12 (37%)	0
26	hM	9/10 (90%)	6 (66%)	0
27	hR	26/27 (96%)	9 (34%)	0
28	hS	16/17 (94%)	14 (87%)	0
All	All	276/284 (97%)	111 (40%)	0

All (111) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
21	bN	63	A
21	bN	65	A
21	bN	73	U
21	bN	74	G
21	bN	86	U
21	bN	88	C
21	bN	89	A
21	bN	91	U
21	bN	95	U
21	bN	96	A
21	bN	98	U
21	bN	99	C
21	bN	100	C
21	bN	102	C
21	bN	103	G
21	bN	104	C
21	bN	105	U
21	bN	106	C
21	bN	107	A
21	bN	110	C
21	bN	119	C
22	hB	11	A
22	hB	15	C
22	hB	18	A
22	hB	22	G
22	hB	23	U
22	hB	25	A
22	hB	26	U
22	hB	27	U
22	hB	35	A

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Mol	Chain	Res	Type
22	hB	38	U
22	hB	41	U
22	hB	43	C
23	hD	2	U
23	hD	3	U
23	hD	4	U
23	hD	7	U
23	hD	8	U
23	hD	12	U
24	hG	6	U
24	hG	7	C
24	hG	8	A
24	hG	9	C
24	hG	11	U
24	hG	12	A
24	hG	14	C
24	hG	16	A
24	hG	17	A
24	hG	28	U
24	hG	29	U
24	hG	30	G
24	hG	38	A
24	hG	39	C
24	hG	40	A
24	hG	41	A
24	hG	42	U
24	hG	44	C
24	hG	45	C
24	hG	46	C
24	hG	47	U
24	hG	48	U
24	hG	49	C
24	hG	50	U
24	hG	51	A
24	hG	52	G
24	hG	55	G
24	hG	57	G
24	hG	59	A
24	hG	60	A
24	hG	72	C
25	hL	10	G
25	hL	12	A

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Mol	Chain	Res	Type
25	hL	18	A
25	hL	19	G
25	hL	20	A
25	hL	22	G
25	hL	23	A
25	hL	24	A
25	hL	25	U
25	hL	26	G
25	hL	27	C
25	hL	30	A
26	hM	2	U
26	hM	3	U
26	hM	4	U
26	hM	7	U
26	hM	8	U
26	hM	10	U
27	hR	10	G
27	hR	13	U
27	hR	14	U
27	hR	15	A
27	hR	18	A
27	hR	19	U
27	hR	20	G
27	hR	21	G
27	hR	31	C
28	hS	2	U
28	hS	3	U
28	hS	4	U
28	hS	5	U
28	hS	8	U
28	hS	9	U
28	hS	10	U
28	hS	11	U
28	hS	12	U
28	hS	13	U
28	hS	14	U
28	hS	15	U
28	hS	16	U
28	hS	17	U

There are no RNA pucker outliers to report.

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 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
24	hG	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	hG	40:A	O3'	41:A	P	1.80

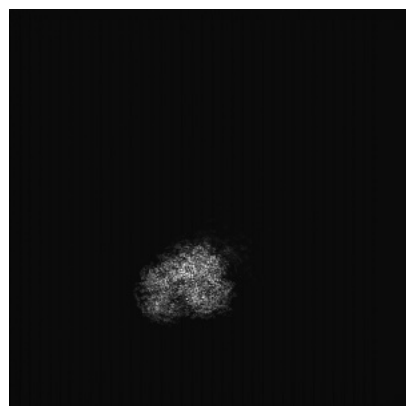
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-50448. These allow visual inspection of the internal detail of the map and identification of artifacts.

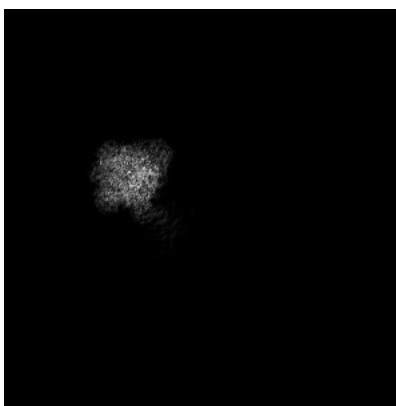
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

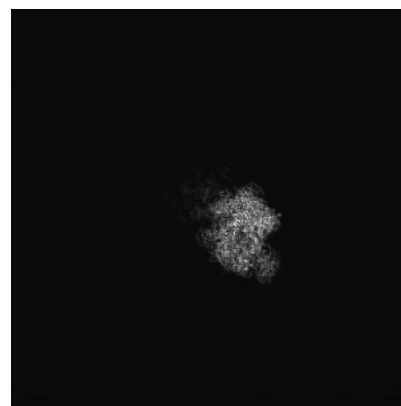
6.1.1 Primary map



X

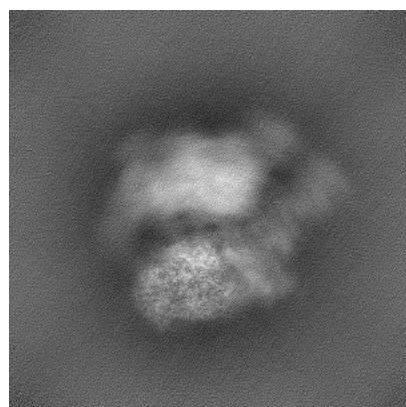


Y

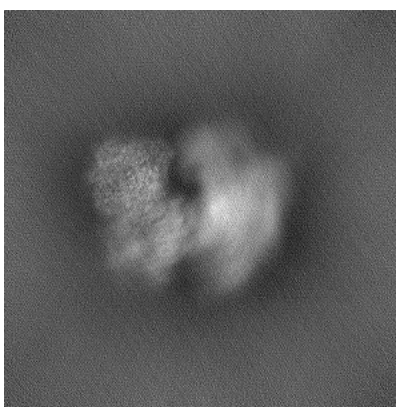


Z

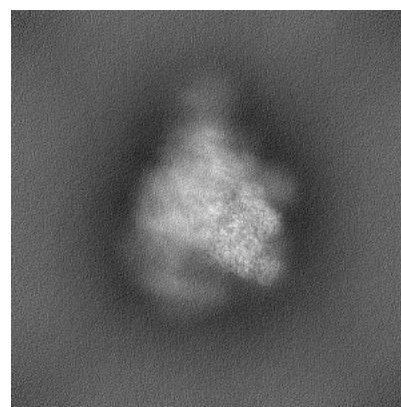
6.1.2 Raw map



X



Y



Z

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

6.2 Central slices [i](#)

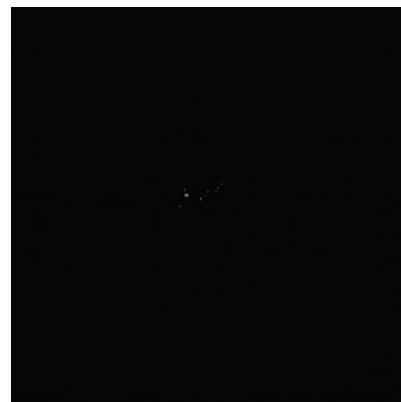
6.2.1 Primary map



X Index: 210



Y Index: 210

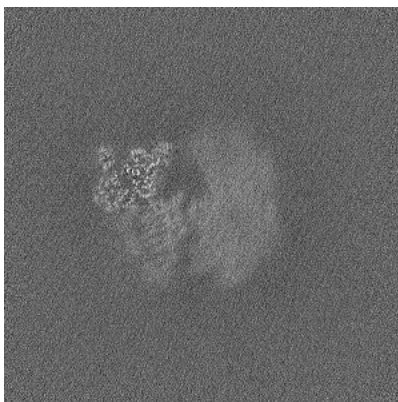


Z Index: 210

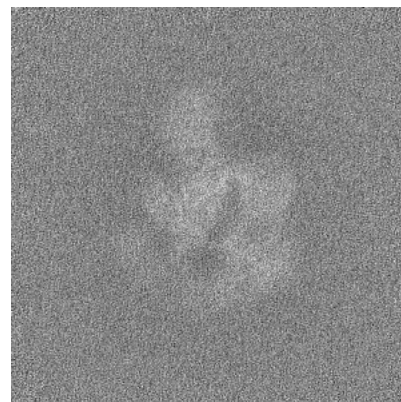
6.2.2 Raw map



X Index: 210



Y Index: 210

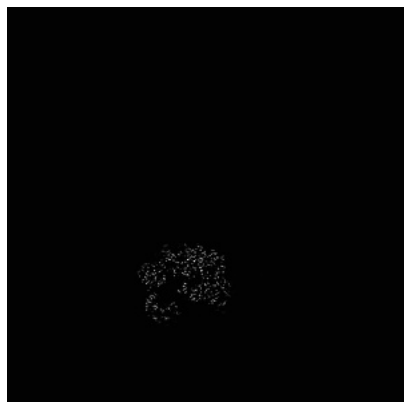


Z Index: 210

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

6.3 Largest variance slices [i](#)

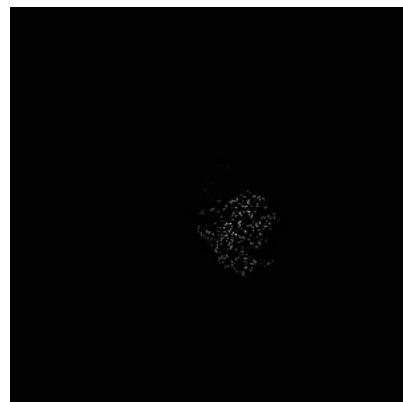
6.3.1 Primary map



X Index: 246

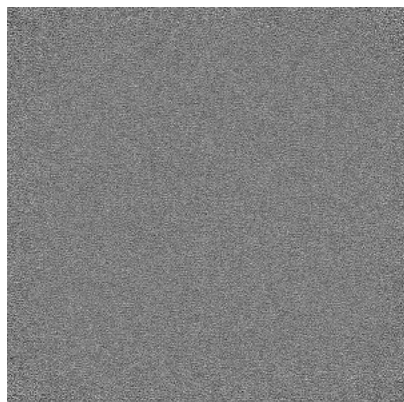


Y Index: 194

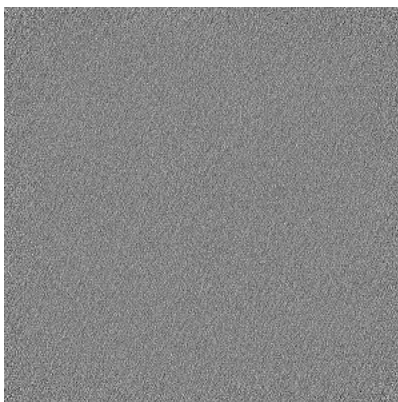


Z Index: 140

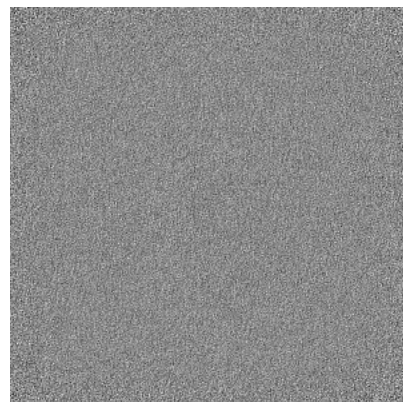
6.3.2 Raw map



X Index: 0



Y Index: 0

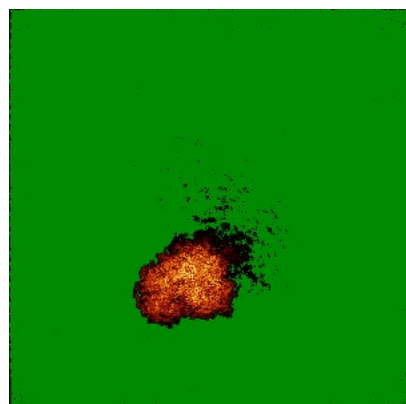


Z Index: 419

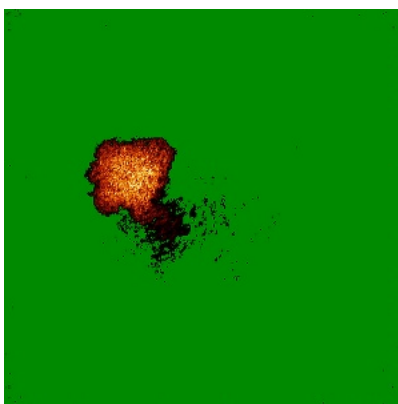
The images above show the largest variance slices of the map in three orthogonal directions.

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

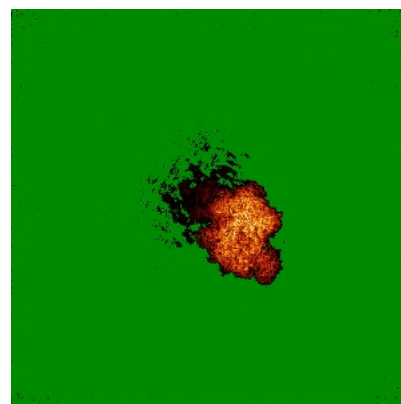
6.4.1 Primary map



X

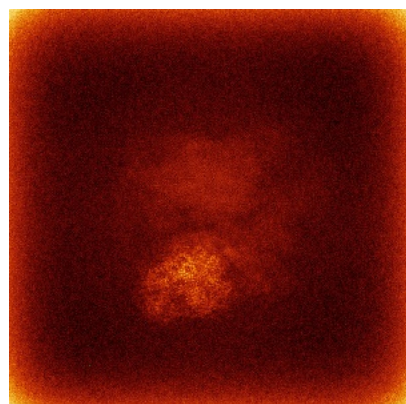


Y

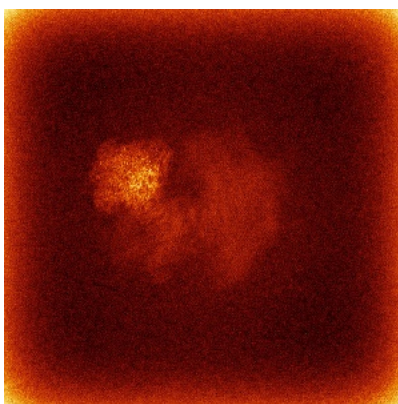


Z

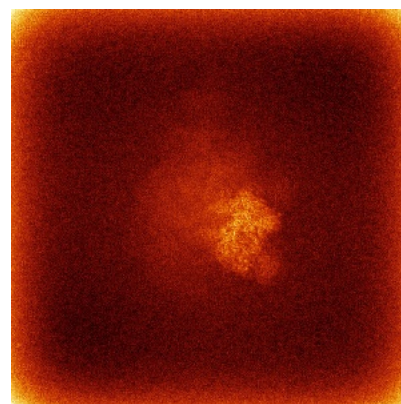
6.4.2 Raw map



X



Y

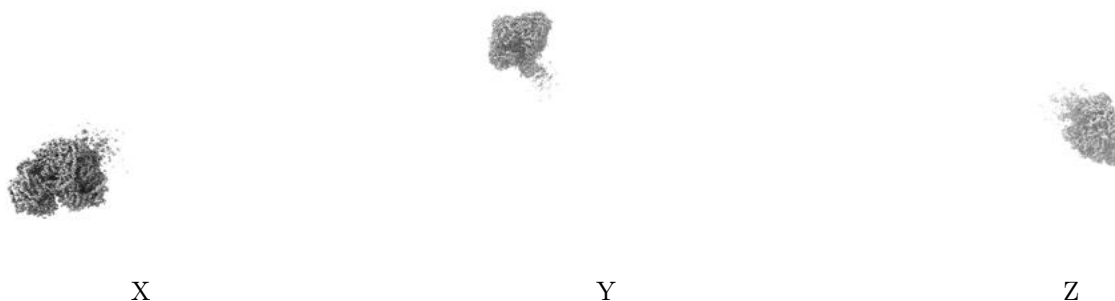


Z

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

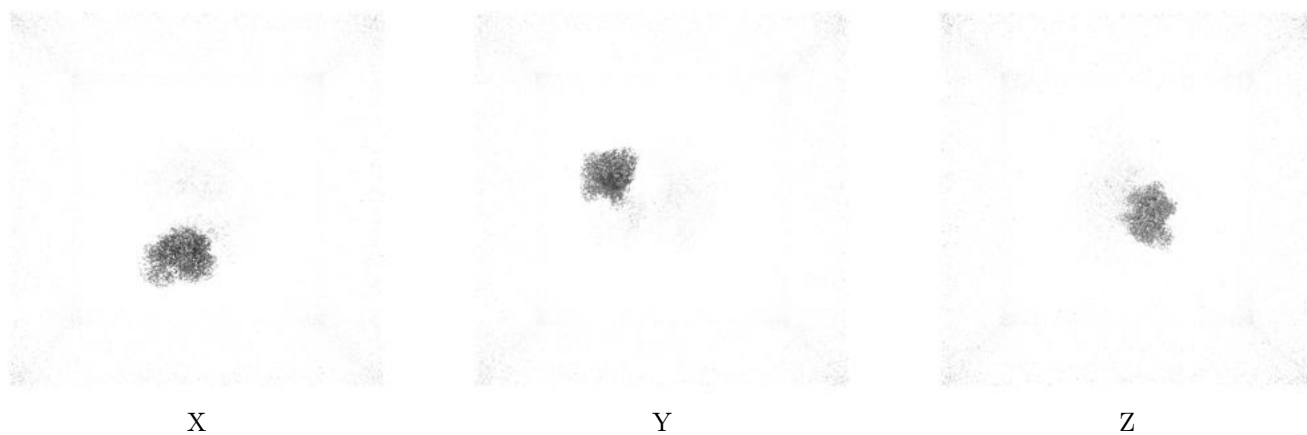
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

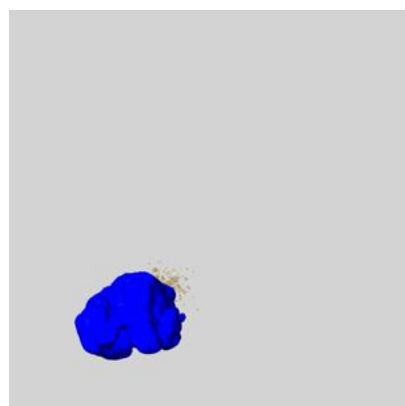
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

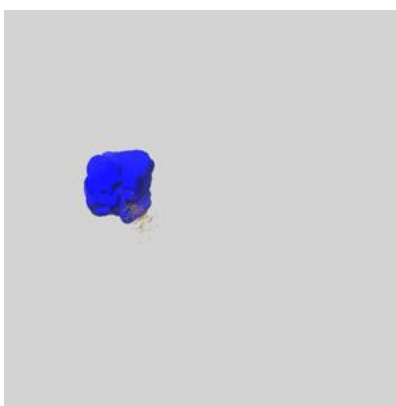
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

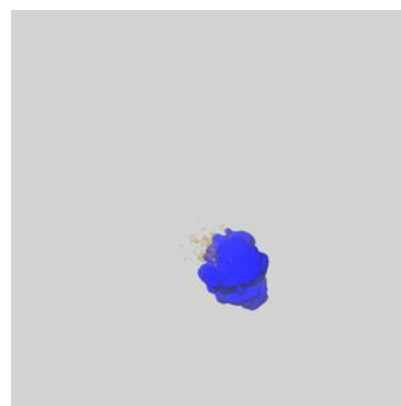
6.6.1 emd_50448_msk_1.map [i](#)



X



Y

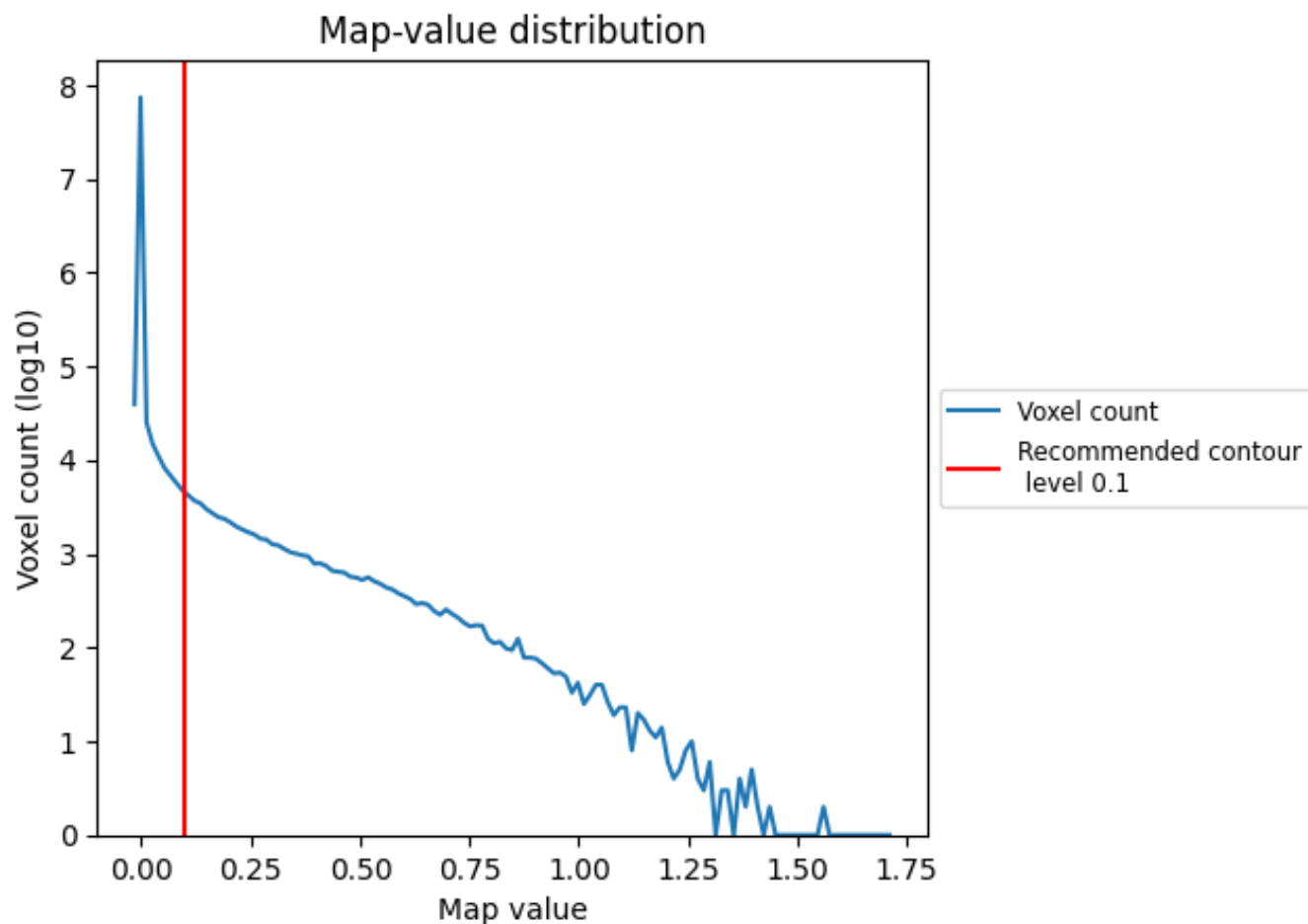


Z

7 Map analysis [i](#)

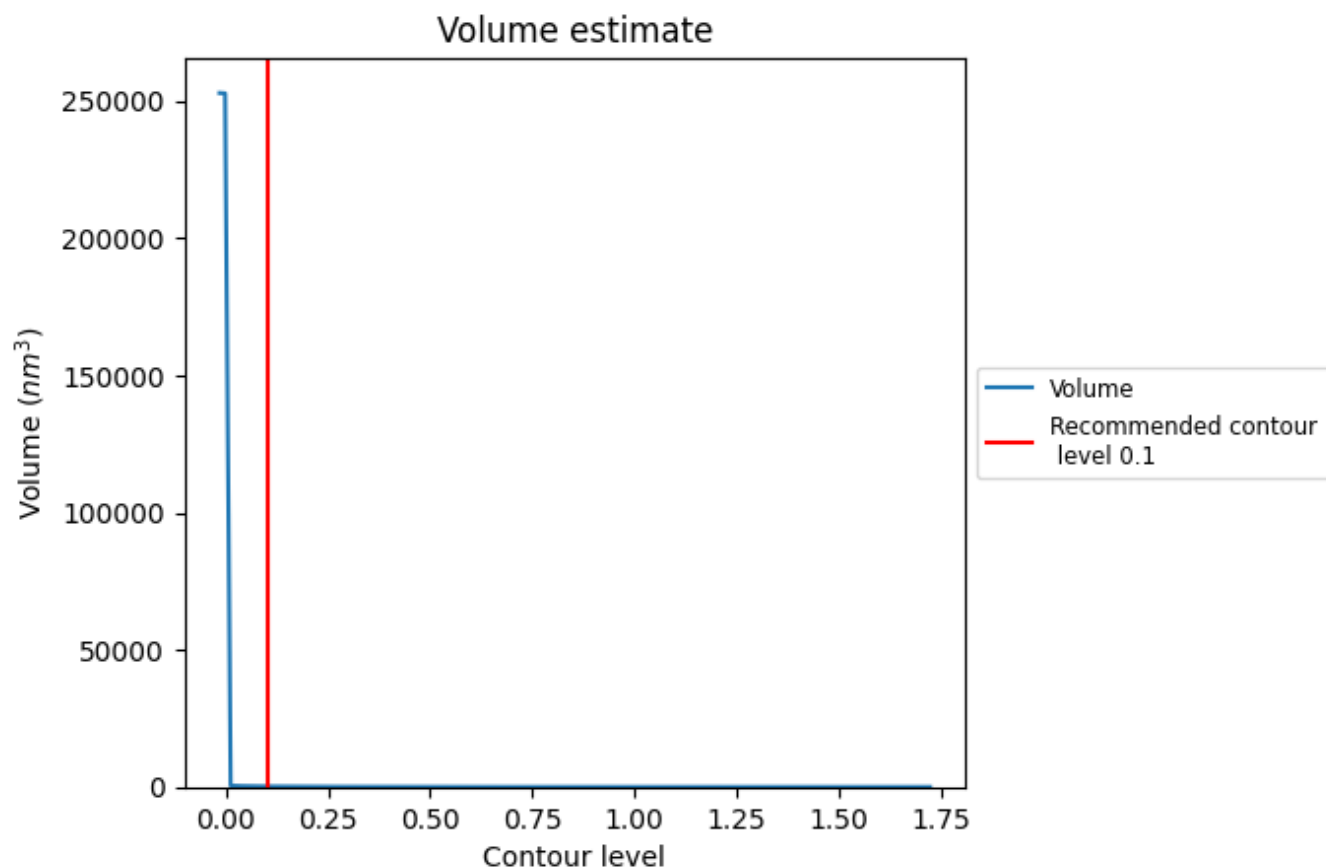
This section contains the results of statistical analysis of the map.

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.

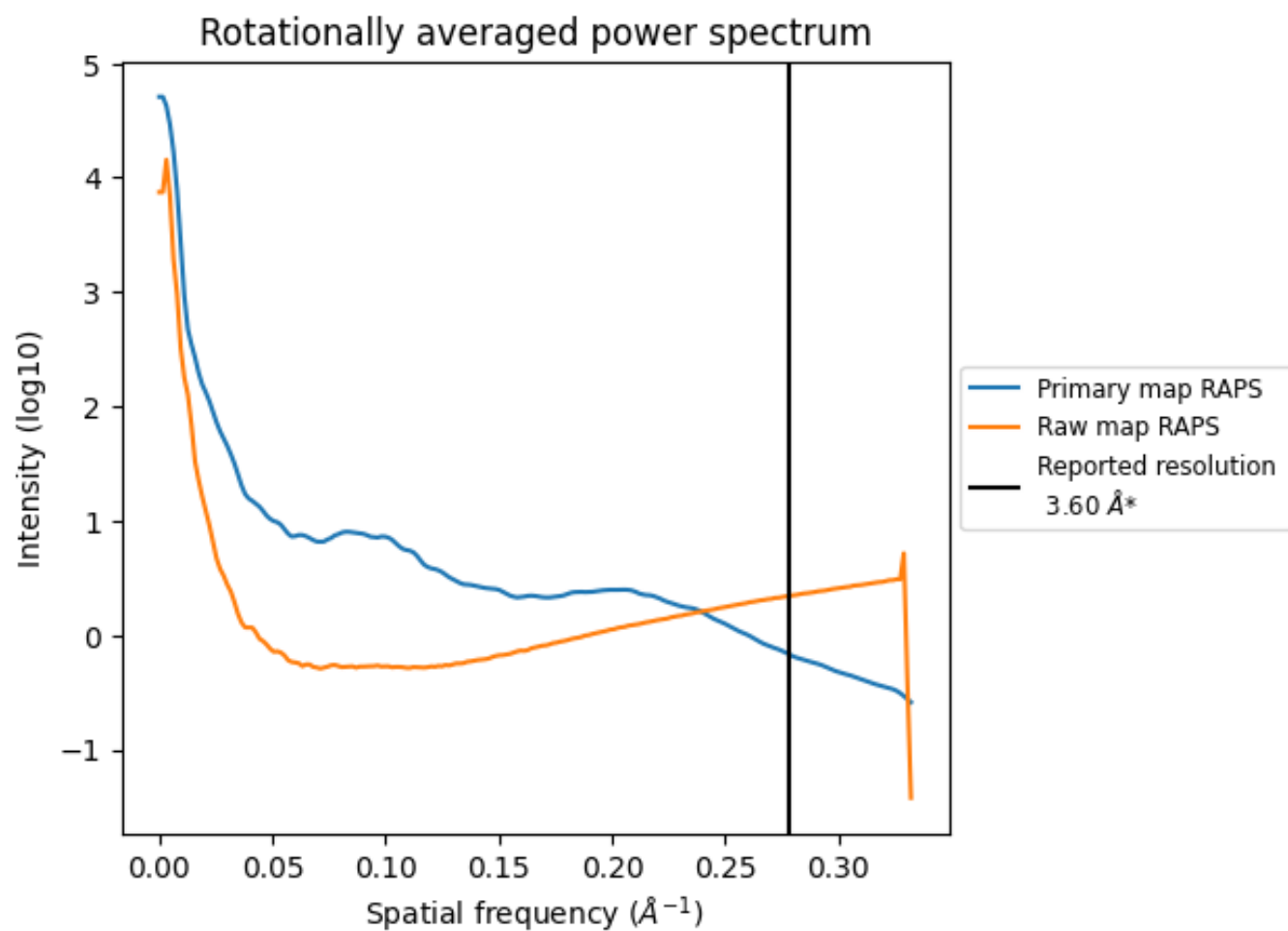
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 198 nm³; this corresponds to an approximate mass of 178 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 ⓘ

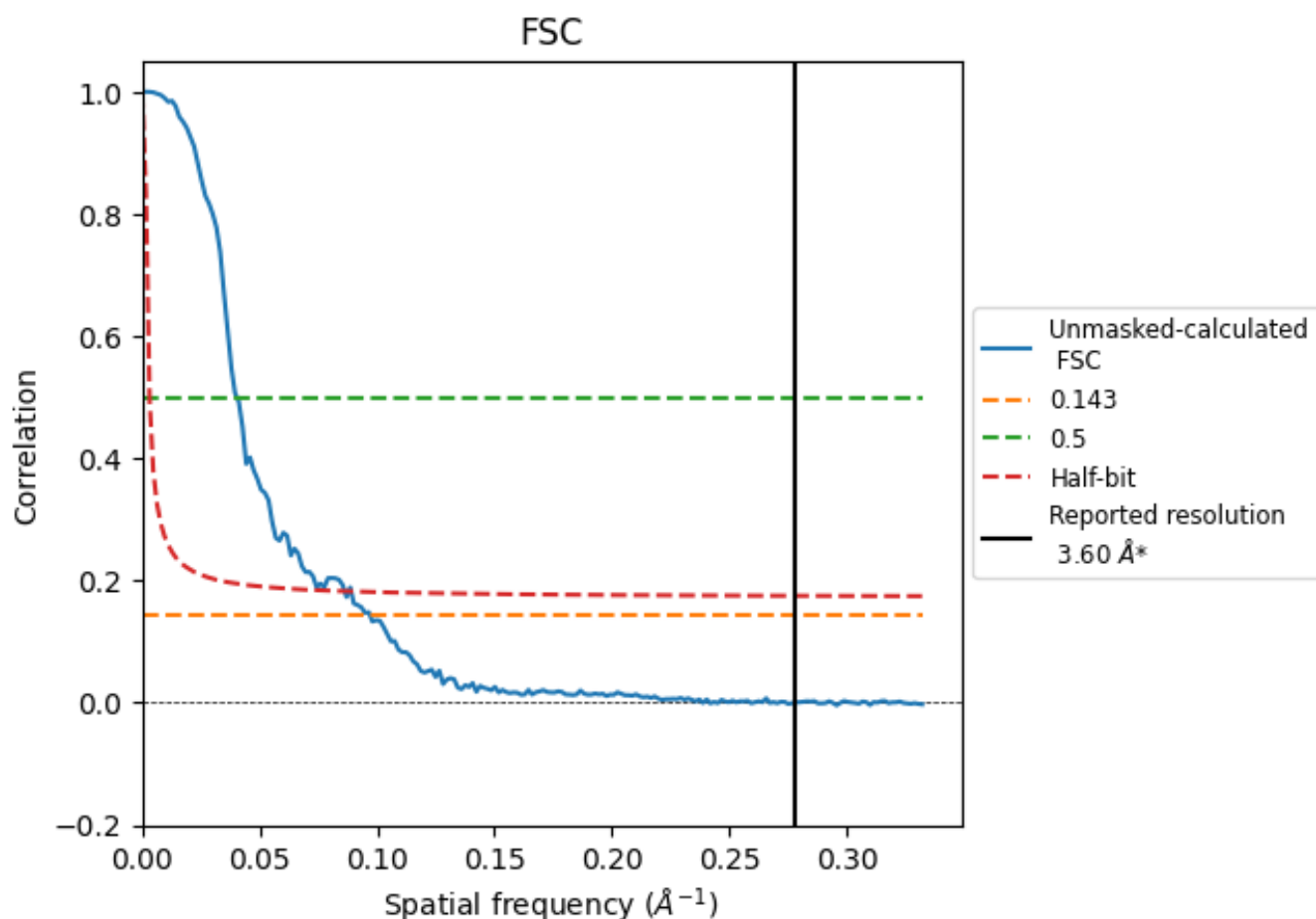


*Reported resolution corresponds to spatial frequency of 0.278 $\text{\AA}^{-1}$

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 $0.278 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

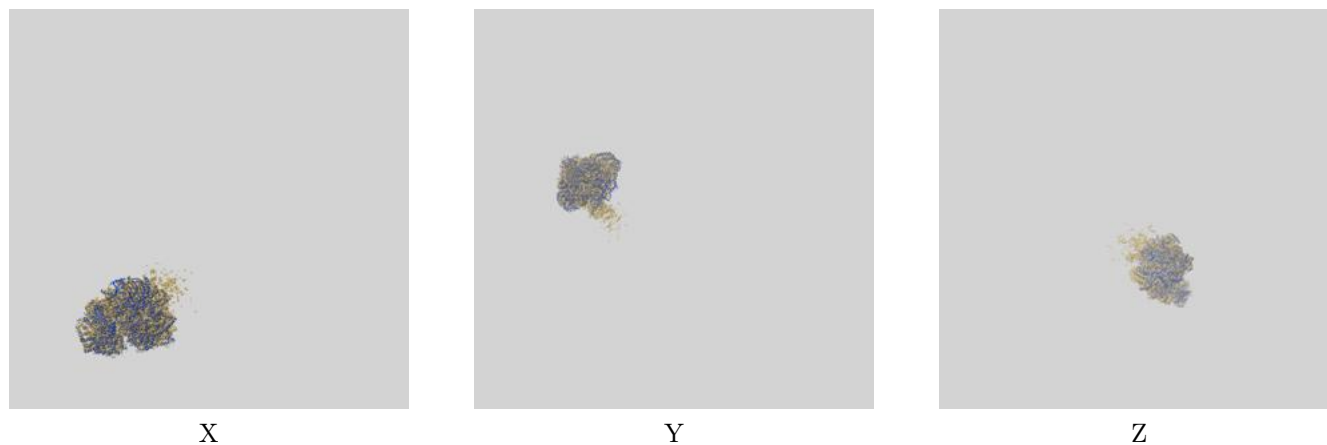
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.60	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	10.31	24.81	11.60

*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 10.31 differs from the reported value 3.6 by more than 10 %

9 Map-model fit [i](#)

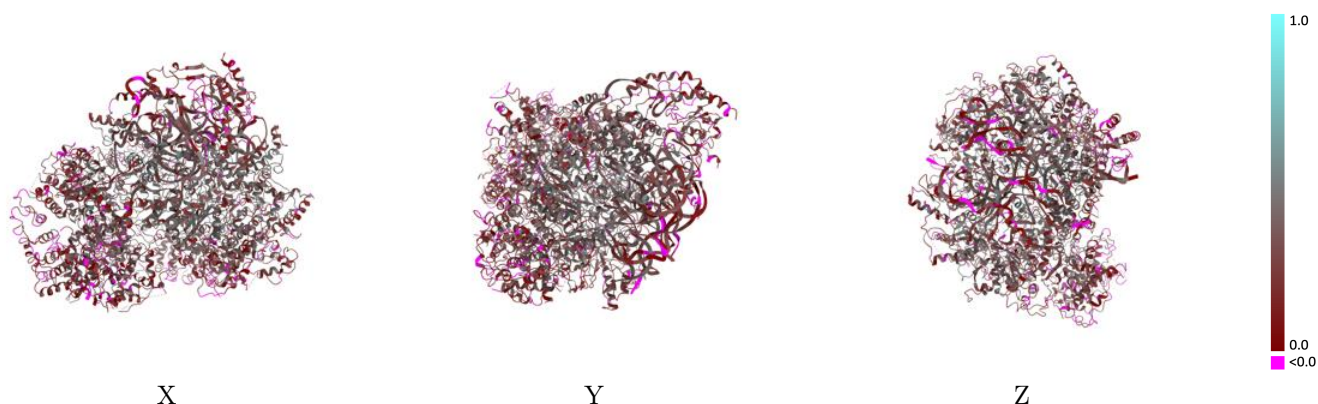
This section contains information regarding the fit between EMDB map EMD-50448 and PDB model 9FI8. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



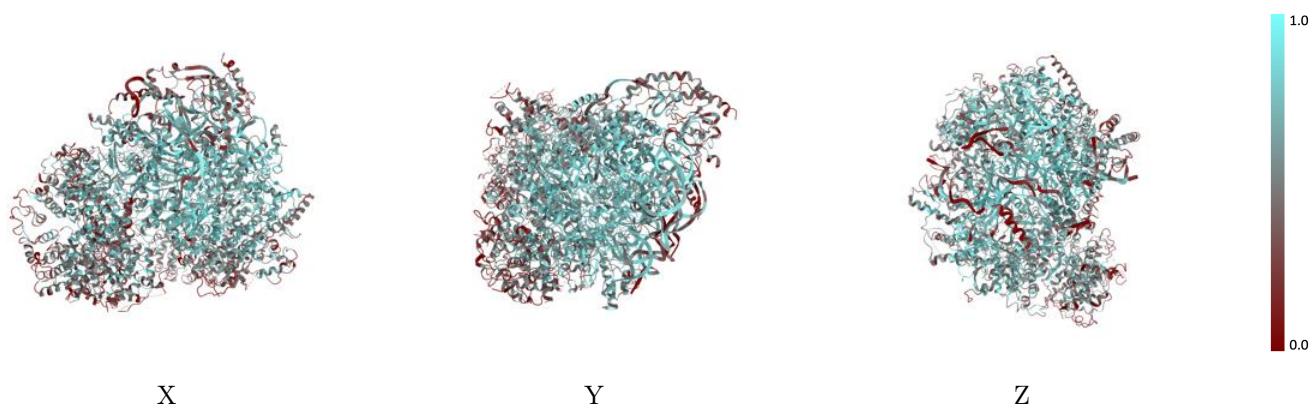
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



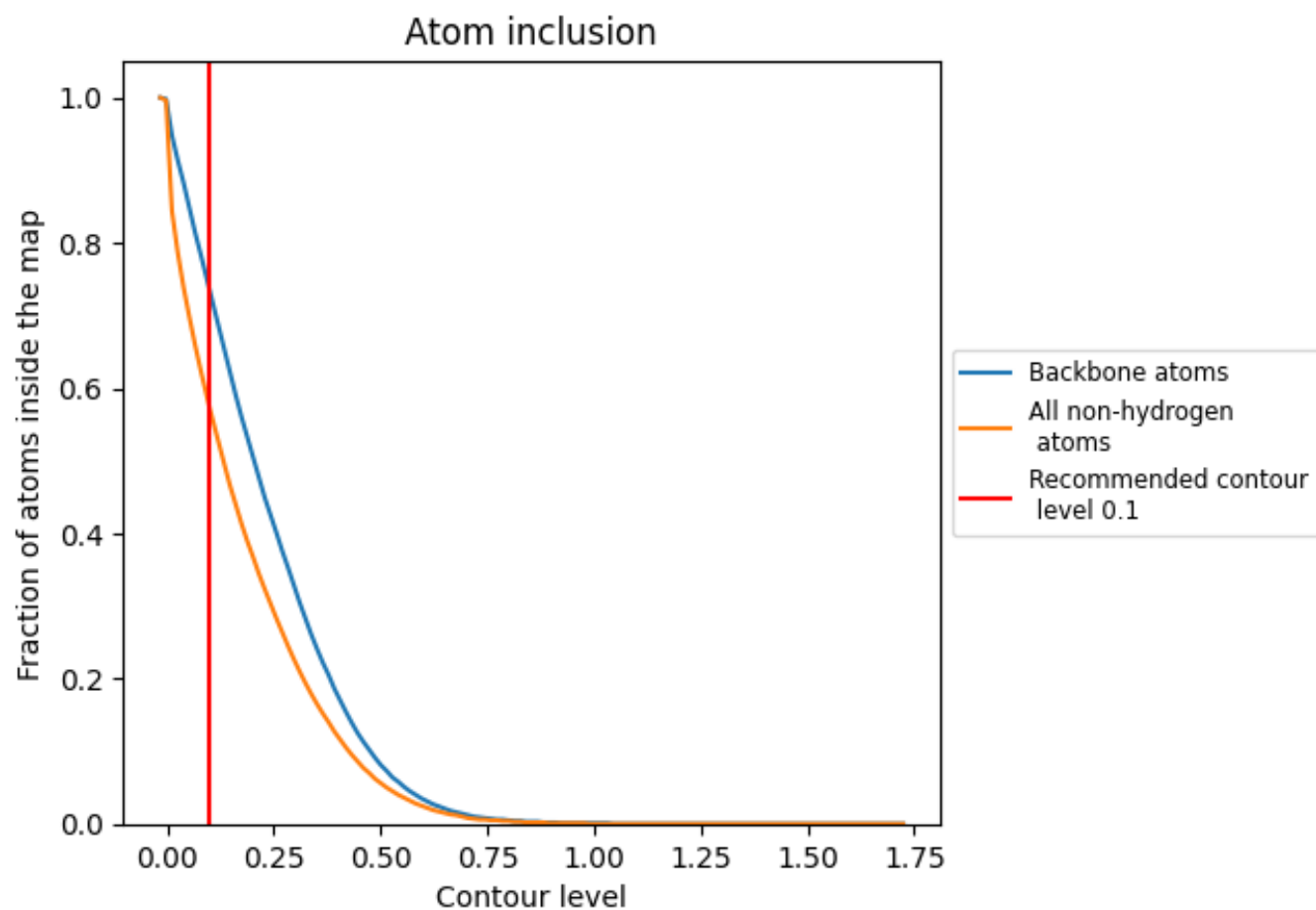
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 57% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5720	 0.2690
HA	 0.6040	 0.2830
HB	 0.6710	 0.3120
HC	 0.6930	 0.3400
HD	 0.8030	 0.4110
HE	 0.6810	 0.3630
HF	 0.6440	 0.3120
HG	 0.4980	 0.2220
HH	 0.6660	 0.3110
HI	 0.7090	 0.3550
HJ	 0.4350	 0.2240
HK	 0.4200	 0.1480
HL	 0.4820	 0.2770
HM	 0.4450	 0.2070
HN	 0.7610	 0.4160
HO	 0.4930	 0.2350
HP	 0.7150	 0.3600
HQ	 0.5020	 0.2320
HR	 0.6130	 0.2680
HS	 0.5810	 0.2800
HT	 0.5350	 0.2700
bN	 0.5060	 0.1980
hB	 0.6600	 0.2690
hD	 0.7430	 0.2820
hG	 0.5620	 0.2160
hL	 0.7650	 0.3310
hM	 0.7450	 0.2760
hR	 0.7600	 0.3200
hS	 0.5340	 0.2940

