



# wwPDB EM Validation Summary Report ⓘ

Jun 29, 2026 – 02:26 pm BST

PDB ID : 9SDP / pdb\_00009sdp  
EMDB ID : EMD-54789  
Title : Cryo-EM structure of the Arabidopsis thaliana 40S ribosomal subunit  
Authors : Karki, S.; Lu, X.; Paatero, A.O.; Ruonala, R.; Tranter, D.; Guryanov, S.;  
Rehan, S.; Hellmann, E.; Haakonsson, A.; Butcher, S.J.; Huiskonen, J.T.;  
Kajander, T.; Helariutta, Y.; Paavilainen, V.O.  
Deposited on : 2025-08-14  
Resolution : 3.30 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

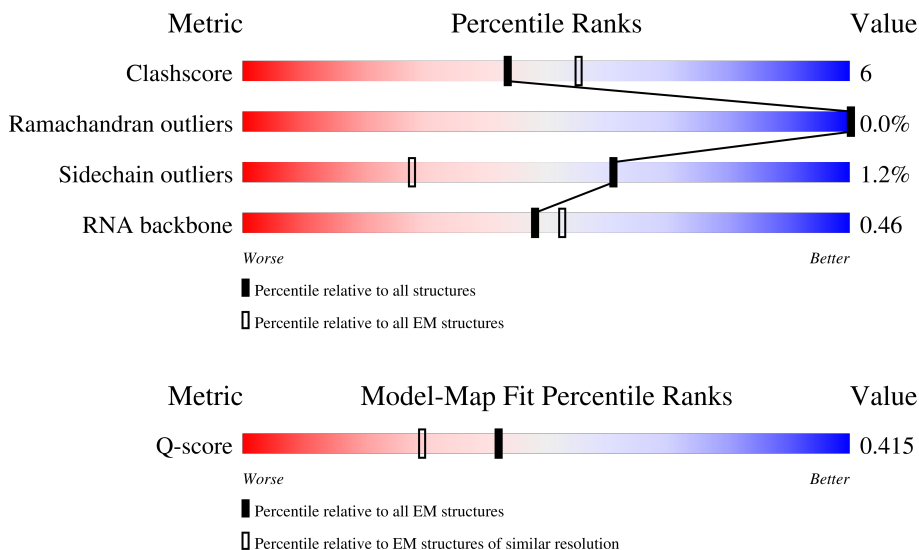
EMDB validation analysis : 0.0.1.dev133  
Mogul : 1.8.4, CSD as541be (2020)  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.50

# 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 3.30 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	15087 ( 2.80 - 3.80 )

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  $\geq 3$ , 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	9	1804	
2	BA	280	
3	BB	262	

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Mol	Chain	Length	Quality of chain
4	BC	276	64% 13% 22%
5	BE	261	80% 16% ..
6	BG	249	12% 66% 14% 20%
7	BH	191	19% 83% 8% 8%
8	BI	222	68% 14% 18%
9	BJ	197	73% 14% 13%
10	BL	160	69% 17% 14%
11	BN	151	87% 12% .
12	cS	150	74% 10% 16%
13	BV	82	72% 16% 12%
14	BW	130	84% 15% .
15	BX	142	85% 12% ..
16	BY	133	63% 19% 17%
17	BZ	130	65% 8% 25%
18	Bb	86	88% 10% .
19	Be	62	56% 6% 37%
20	Cn	25	12% 76% 12% 12%

## 2 Entry composition [i](#)

There are 23 unique types of molecules in this entry. The entry contains 44802 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 18S RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	9	1081	23135	10353	4182	7519	1081	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
9	604	UY1	U	modified residue	GB X16077.1

- Molecule 2 is a protein called Small ribosomal subunit protein uS2y.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	BA	189	1500	957	263	268	12	0	0

- Molecule 3 is a protein called Small ribosomal subunit protein eS1z.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	BB	208	1696	1075	312	301	8	0	0

- Molecule 4 is a protein called Small ribosomal subunit protein uS5w.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	BC	216	1683	1088	298	290	7	0	0

- Molecule 5 is a protein called Small ribosomal subunit protein eS4z.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	BE	255	2058	1316	382	354	6	0	0

- Molecule 6 is a protein called Small ribosomal subunit protein eS6y.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	BG	199	1579	993	310	268	8	0	0

- Molecule 7 is a protein called Small ribosomal subunit protein eS7z.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	BH	176	1439	918	262	258	1	0	0

- Molecule 8 is a protein called Small ribosomal subunit protein eS8z.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	BI	183	1480	921	293	262	4	0	0

- Molecule 9 is a protein called Small ribosomal subunit protein uS4y.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	BJ	171	1432	906	280	242	4	0	0

- Molecule 10 is a protein called Small ribosomal subunit protein uS17z.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	BL	138	1106	709	214	178	5	0	0

- Molecule 11 is a protein called Small ribosomal subunit protein uS15y.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	BN	149	1192	760	223	207	2	0	0

- Molecule 12 is a protein called Small ribosomal subunit protein uS11x.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	cS	126	952	586	188	174	4	0	0

- Molecule 13 is a protein called Small ribosomal subunit protein eS21z.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	BV	72	Total	C	N	O	S	0	0
			561	350	103	105	3		

- Molecule 14 is a protein called Small ribosomal subunit protein uS8z/uS8w.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	BW	129	Total	C	N	O	S	0	0
			1033	660	188	180	5		

- Molecule 15 is a protein called Small ribosomal subunit protein uS12y.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	BX	138	Total	C	N	O	S	0	0
			1075	684	206	182	3		

- Molecule 16 is a protein called Small ribosomal subunit protein eS24z.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	BY	110	Total	C	N	O	S	0	0
			917	587	176	152	2		

- Molecule 17 is a protein called Small ribosomal subunit protein eS26x.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	BZ	97	Total	C	N	O	S	0	0
			783	482	165	130	6		

- Molecule 18 is a protein called Small ribosomal subunit protein eS27y.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Bb	85	Total	C	N	O	S	0	0
			660	411	120	122	7		

- Molecule 19 is a protein called Small ribosomal subunit protein eS30z/eS30y/eS30x.

Mol	Chain	Residues	Atoms				AltConf	Trace
19	Be	39	Total	C	N	O	0	0
			292	175	70	47		

- Molecule 20 is a protein called Small ribosomal subunit protein eS32 eS32z/eS32y/eS32x/eS32w/eS32v.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
20	Cn	22	211	130	55	23	3	0	0

- Molecule 21 is POTASSIUM ION (CCD ID: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
21	9	3	Total	K	0
			3	3	

- Molecule 22 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
22	9	14	Total	Mg	0
			14	14	

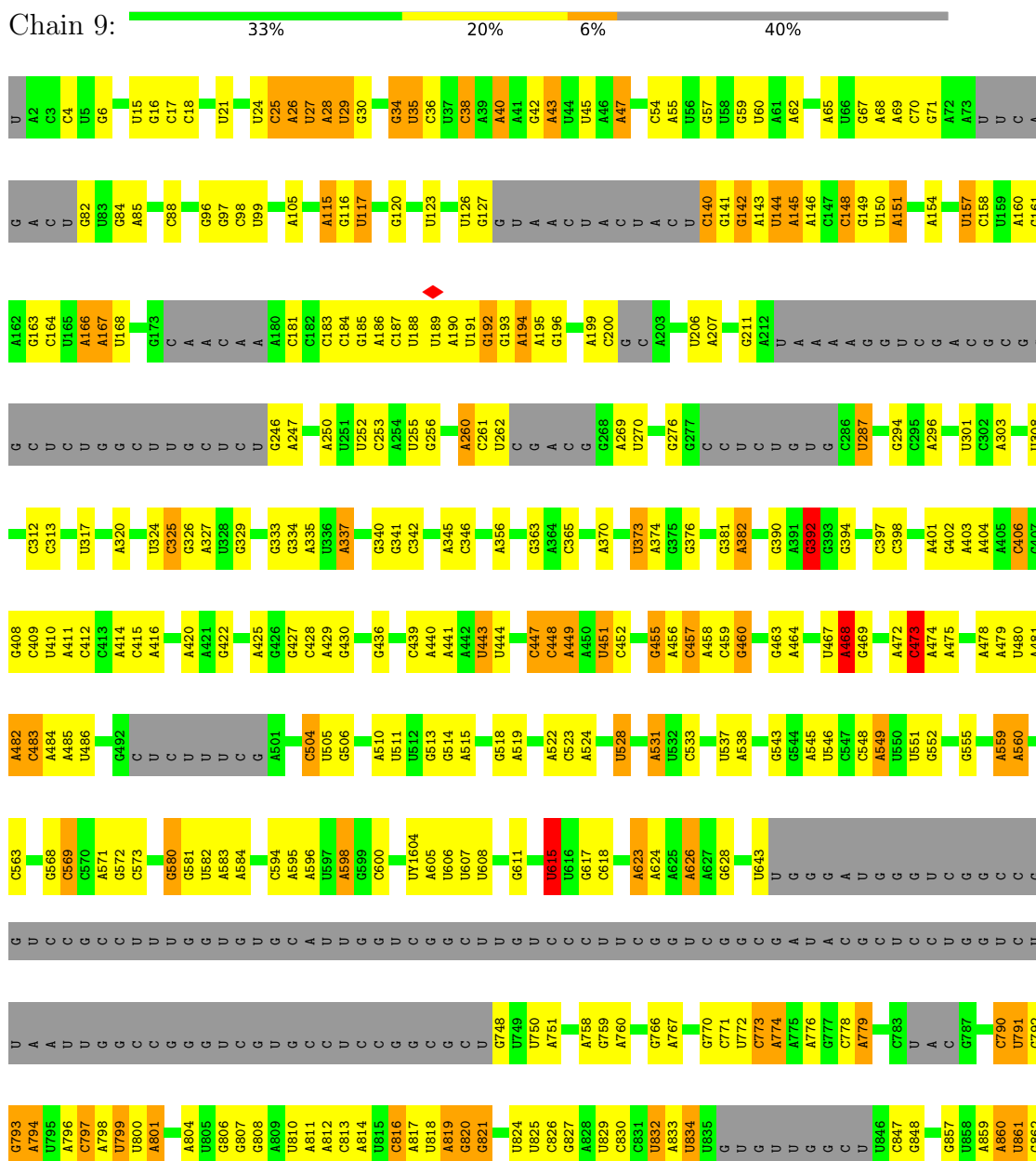
- Molecule 23 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

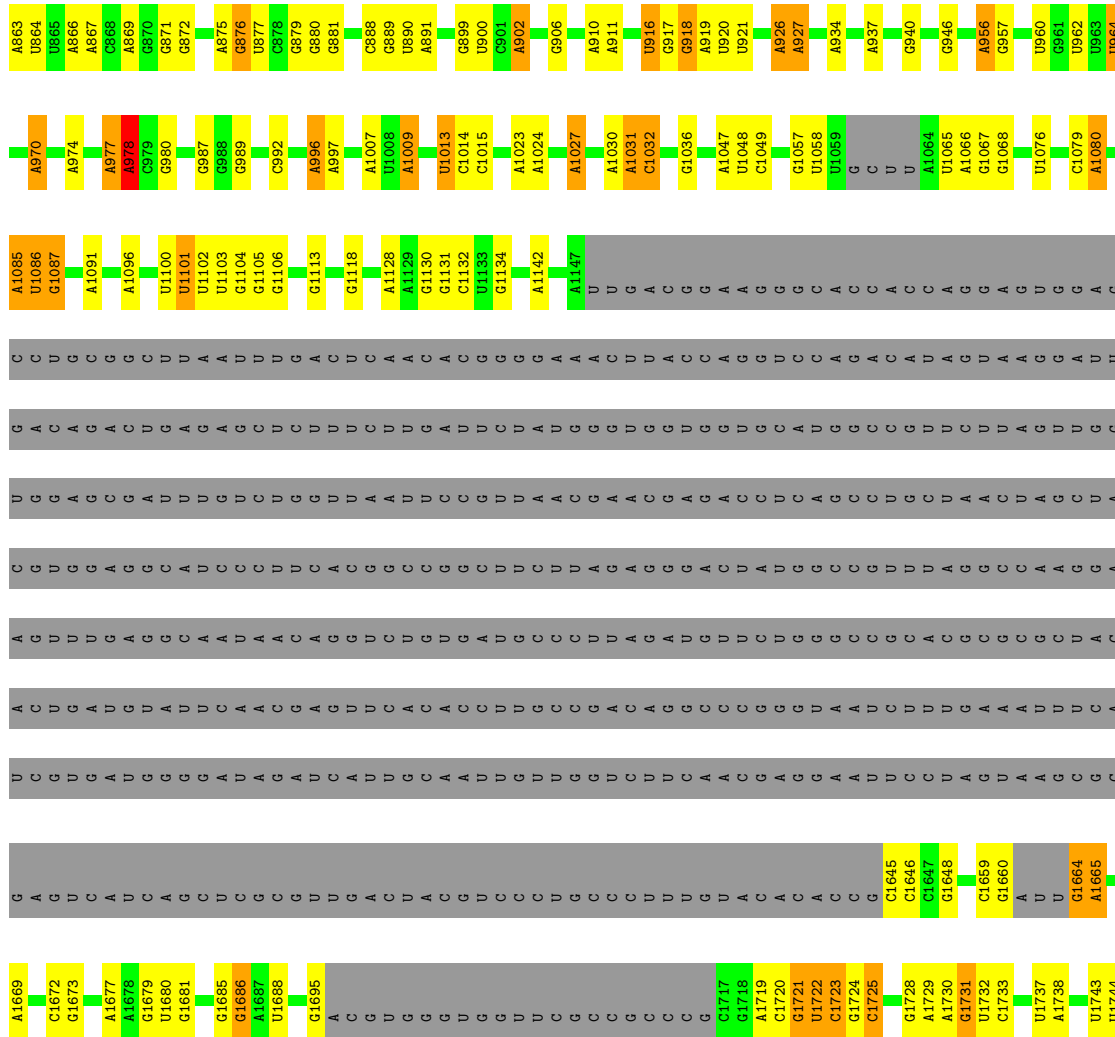
Mol	Chain	Residues	Atoms		AltConf
23	BZ	1	Total	Zn	0
			1	1	

### 3 Residue-property plots

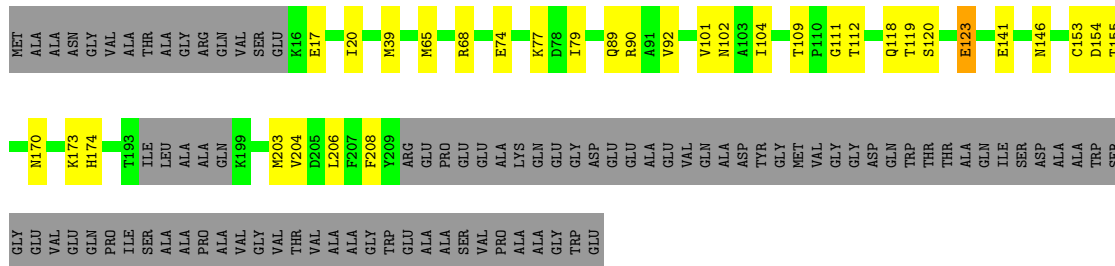
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: 18S RNA





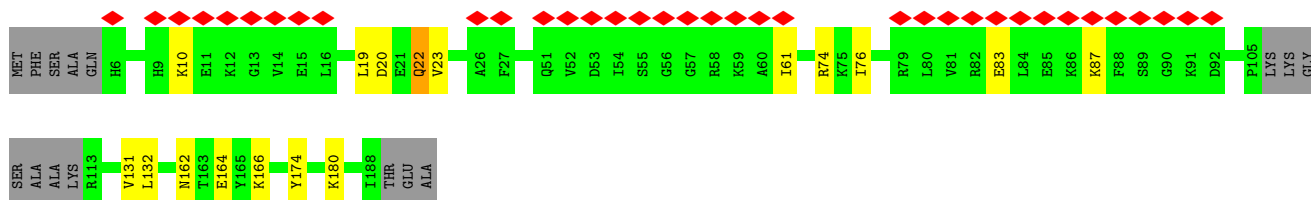
● Molecule 2: Small ribosomal subunit protein uS2y



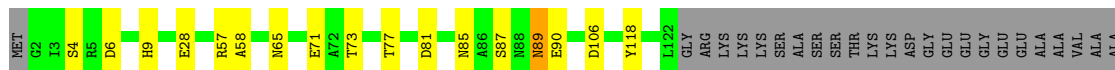
● Molecule 3: Small ribosomal subunit protein eS1z



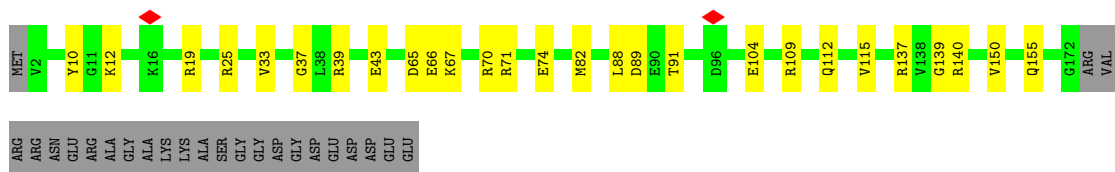
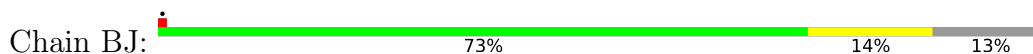




• Molecule 8: Small ribosomal subunit protein eS8z



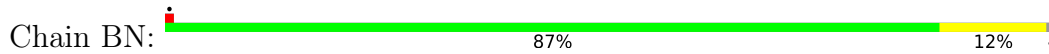
• Molecule 9: Small ribosomal subunit protein uS4y



• Molecule 10: Small ribosomal subunit protein uS17z



• Molecule 11: Small ribosomal subunit protein uS15y

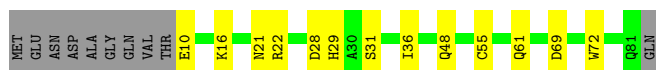


• Molecule 12: Small ribosomal subunit protein uS11x

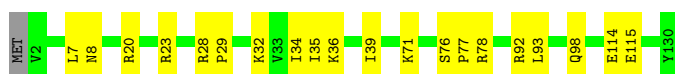
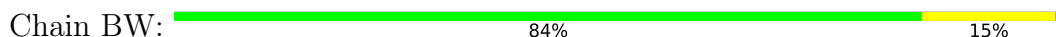




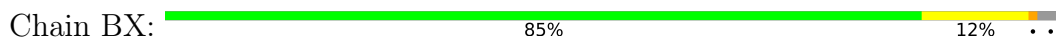
- Molecule 13: Small ribosomal subunit protein eS21z



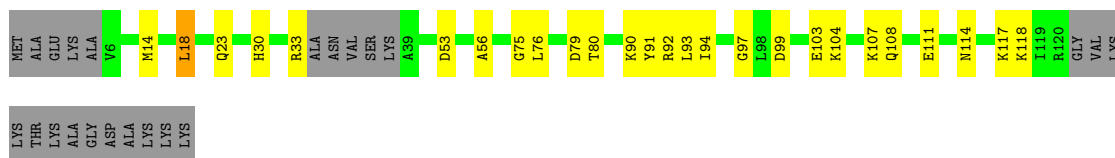
- Molecule 14: Small ribosomal subunit protein uS8z/uS8w



- Molecule 15: Small ribosomal subunit protein uS12y



- Molecule 16: Small ribosomal subunit protein eS24z



- Molecule 17: Small ribosomal subunit protein eS26x

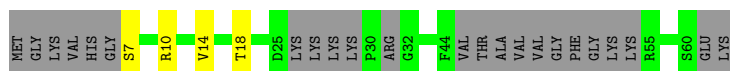


- Molecule 18: Small ribosomal subunit protein eS27y




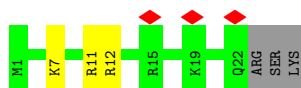
- Molecule 19: Small ribosomal subunit protein eS30z/eS30y/eS30x

Chain Be:  56% 6% 37%



- Molecule 20: Small ribosomal subunit protein eS32 eS32z/eS32y/eS32x/eS32w/eS32v

Chain Cn:  12% 76% 12% 12%



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	115070	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	24.512	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1200	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.319	Depositor
Minimum map value	-0.116	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.017	Depositor
Recommended contour level	0.06	Depositor
Map size ( $\text{\AA}$ )	378.0, 378.0, 378.0	wwPDB
Map dimensions	450, 450, 450	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.84, 0.84, 0.84	Depositor

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: MG, MA6, 4AC, A2M, ZN, OMC, UY1, OMU, OMG, K

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	9	0.22	0/25437	0.29	0/39618
2	BA	0.19	0/1532	0.33	0/2070
3	BB	0.17	0/1724	0.31	0/2313
4	BC	0.21	0/1719	0.35	0/2319
5	BE	0.17	0/2100	0.32	0/2819
6	BG	0.16	0/1599	0.31	0/2127
7	BH	0.15	0/1461	0.27	0/1964
8	BI	0.16	0/1503	0.31	0/2008
9	BJ	0.18	0/1458	0.35	0/1954
10	BL	0.19	0/1132	0.35	0/1514
11	BN	0.18	0/1216	0.28	0/1632
12	cS	0.18	0/964	0.32	0/1293
13	BV	0.17	0/570	0.29	0/766
14	BW	0.24	0/1051	0.33	0/1406
15	BX	0.19	0/1092	0.35	0/1453
16	BY	0.16	0/931	0.31	0/1233
17	BZ	0.38	0/797	0.53	1/1069 (0.1%)
18	Bb	0.18	0/670	0.28	0/899
19	Be	0.14	0/294	0.23	0/387
20	Cn	0.27	0/212	0.47	0/269
All	All	0.21	0/47462	0.30	1/69113 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	BZ	73	TYR	CA-C-O	-5.01	116.18	121.94

There are no chirality outliers.

There are no planarity outliers.

## 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	9	23135	0	11665	270	0
2	BA	1500	0	1507	21	0
3	BB	1696	0	1760	15	0
4	BC	1683	0	1782	28	0
5	BE	2058	0	2163	39	0
6	BG	1579	0	1694	22	0
7	BH	1439	0	1508	11	0
8	BI	1480	0	1527	23	0
9	BJ	1432	0	1485	17	0
10	BL	1106	0	1169	21	0
11	BN	1192	0	1276	13	0
12	cS	952	0	988	11	0
13	BV	561	0	557	10	0
14	BW	1033	0	1070	16	0
15	BX	1075	0	1146	16	0
16	BY	917	0	981	17	0
17	BZ	783	0	810	10	0
18	Bb	660	0	672	8	0
19	Be	292	0	288	2	0
20	Cn	211	0	258	2	0
21	9	3	0	0	0	0
22	9	14	0	0	0	0
23	BZ	1	0	0	0	0
All	All	44802	0	34306	501	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

The worst 5 of 501 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)
5:BE:165:GLU:OE1	5:BE:166:GLU:HG3	1.50	1.08
1:9:276:G:H1	1:9:287:U:H3	1.16	0.93
1:9:1672:C:N3	1:9:1743:U:O4	2.02	0.92
9:BJ:39:ARG:NH1	9:BJ:43:GLU:OE2	2.03	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:9:1105:G:O4'	14:BW:8:ASN:ND2	2.06	0.89

There are no symmetry-related clashes.

## 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	
2	BA	185/280 (66%)	181 (98%)	4 (2%)	0	100	100
3	BB	206/262 (79%)	194 (94%)	12 (6%)	0	100	100
4	BC	214/276 (78%)	204 (95%)	10 (5%)	0	100	100
5	BE	253/261 (97%)	236 (93%)	17 (7%)	0	100	100
6	BG	191/249 (77%)	182 (95%)	8 (4%)	1 (0%)	24	55
7	BH	172/191 (90%)	170 (99%)	2 (1%)	0	100	100
8	BI	179/222 (81%)	175 (98%)	4 (2%)	0	100	100
9	BJ	169/197 (86%)	160 (95%)	9 (5%)	0	100	100
10	BL	136/160 (85%)	133 (98%)	3 (2%)	0	100	100
11	BN	147/151 (97%)	144 (98%)	3 (2%)	0	100	100
12	cS	124/150 (83%)	116 (94%)	8 (6%)	0	100	100
13	BV	70/82 (85%)	70 (100%)	0	0	100	100
14	BW	127/130 (98%)	122 (96%)	5 (4%)	0	100	100
15	BX	136/142 (96%)	130 (96%)	6 (4%)	0	100	100
16	BY	106/133 (80%)	103 (97%)	3 (3%)	0	100	100
17	BZ	95/130 (73%)	91 (96%)	4 (4%)	0	100	100
18	Bb	83/86 (96%)	82 (99%)	1 (1%)	0	100	100
19	Be	32/62 (52%)	32 (100%)	0	0	100	100
20	Cn	20/25 (80%)	19 (95%)	1 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2645/3189 (83%)	2544 (96%)	100 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	BG	14	LYS

### 5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	BA	161/221 (73%)	159 (99%)	2 (1%)	63	75
3	BB	185/226 (82%)	184 (100%)	1 (0%)	81	83
4	BC	185/225 (82%)	183 (99%)	2 (1%)	65	76
5	BE	224/228 (98%)	218 (97%)	6 (3%)	39	63
6	BG	170/213 (80%)	170 (100%)	0	100	100
7	BH	159/169 (94%)	157 (99%)	2 (1%)	61	74
8	BI	156/183 (85%)	154 (99%)	2 (1%)	61	74
9	BJ	153/172 (89%)	152 (99%)	1 (1%)	76	80
10	BL	119/135 (88%)	118 (99%)	1 (1%)	73	79
11	BN	132/133 (99%)	131 (99%)	1 (1%)	73	79
12	cS	99/120 (82%)	97 (98%)	2 (2%)	48	68
13	BV	59/67 (88%)	58 (98%)	1 (2%)	53	71
14	BW	111/112 (99%)	111 (100%)	0	100	100
15	BX	110/114 (96%)	108 (98%)	2 (2%)	51	70
16	BY	98/114 (86%)	96 (98%)	2 (2%)	48	68
17	BZ	86/108 (80%)	83 (96%)	3 (4%)	32	58
18	Bb	77/78 (99%)	77 (100%)	0	100	100
19	Be	27/49 (55%)	27 (100%)	0	100	100

*Continued on next page...*

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
20	Cn	21/24 (88%)	21 (100%)	0	100	100
All	All	2332/2691 (87%)	2304 (99%)	28 (1%)	61	75

5 of 28 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
8	BI	90	GLU
17	BZ	76	SER
11	BN	117	LEU
16	BY	76	LEU
10	BL	112	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
9	BJ	141	GLN
10	BL	34	GLN
17	BZ	7	ASN
10	BL	99	HIS
5	BE	17	HIS

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	9	1066/1804 (59%)	249 (23%)	13 (1%)

5 of 249 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	9	4	C
1	9	6	G
1	9	25	C
1	9	26	A
1	9	27	U

5 of 13 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	9	820	G

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	9	926	A
1	9	1769	U
1	9	1746	U
1	9	1763	A

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

18 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	A2M	9	468	1	22,25,26	3.39	9 (40%)	31,36,39	2.84	10 (32%)
1	OMG	9	392	1	23,26,27	2.62	9 (39%)	33,38,41	2.03	11 (33%)
1	OMU	9	123	1	19,22,23	2.54	7 (36%)	26,31,34	2.20	5 (19%)
1	OMG	9	246	1	23,26,27	2.69	8 (34%)	33,38,41	2.20	11 (33%)
1	OMC	9	140	1	19,22,23	3.03	8 (42%)	26,31,34	0.61	0
1	OMU	9	615	1	19,22,23	2.67	6 (31%)	26,31,34	2.08	7 (26%)
1	OMU	9	1013	1	19,22,23	2.52	7 (36%)	26,31,34	2.31	8 (30%)
1	A2M	9	28	1	22,25,26	3.41	9 (40%)	31,36,39	2.63	10 (32%)
1	OMC	9	1645	1	19,22,23	2.95	8 (42%)	26,31,34	0.66	0
1	4AC	9	1781	1	21,24,25	3.55	9 (42%)	29,34,37	1.34	4 (13%)
1	UY1	9	604	1	19,22,23	4.09	8 (42%)	22,31,34	2.40	5 (22%)
1	OMC	9	38	1	19,22,23	2.98	8 (42%)	26,31,34	0.64	0
1	A2M	9	978	1	22,25,26	3.43	10 (45%)	31,36,39	2.62	10 (32%)
1	MA6	9	1789	1	23,26,27	1.32	3 (13%)	34,38,41	2.98	13 (38%)
1	A2M	9	1758	1	22,25,26	3.39	9 (40%)	31,36,39	2.60	10 (32%)
1	MA6	9	1790	1	23,26,27	1.31	3 (13%)	34,38,41	3.05	11 (32%)
1	OMC	9	473	1	19,22,23	2.99	8 (42%)	26,31,34	0.78	1 (3%)
1	A2M	9	623	1,22	22,25,26	3.38	10 (45%)	31,36,39	2.91	11 (35%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	A2M	9	468	1	-	2/9/27/28	0/3/3/3
1	OMG	9	392	1	-	3/9/27/28	0/3/3/3
1	OMU	9	123	1	-	0/9/27/28	0/2/2/2
1	OMG	9	246	1	-	2/9/27/28	0/3/3/3
1	OMC	9	140	1	-	3/9/27/28	0/2/2/2
1	OMU	9	615	1	-	5/9/27/28	0/2/2/2
1	OMU	9	1013	1	-	1/9/27/28	0/2/2/2
1	A2M	9	28	1	-	3/9/27/28	0/3/3/3
1	OMC	9	1645	1	-	0/9/27/28	0/2/2/2
1	4AC	9	1781	1	-	2/11/29/30	0/2/2/2
1	UY1	9	604	1	-	2/9/27/28	0/2/2/2
1	OMC	9	38	1	-	2/9/27/28	0/2/2/2
1	A2M	9	978	1	-	2/9/27/28	0/3/3/3
1	MA6	9	1789	1	-	3/11/29/30	0/3/3/3
1	A2M	9	1758	1	-	2/9/27/28	0/3/3/3
1	MA6	9	1790	1	-	6/11/29/30	0/3/3/3
1	OMC	9	473	1	-	3/9/27/28	0/2/2/2
1	A2M	9	623	1,22	-	2/9/27/28	0/3/3/3

The worst 5 of 139 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	9	604	UY1	C6-C5	10.62	1.47	1.35
1	9	604	UY1	C2-N1	9.24	1.49	1.36
1	9	978	A2M	C2'-C1'	-8.53	1.31	1.53
1	9	1758	A2M	C2'-C1'	-8.52	1.31	1.53
1	9	468	A2M	O4'-C1'	8.51	1.62	1.42

The worst 5 of 127 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	9	1790	MA6	N1-C6-N6	-11.52	104.49	117.08
1	9	1789	MA6	N1-C6-N6	-11.38	104.64	117.08
1	9	623	A2M	N6-C6-N1	-8.71	99.27	118.35
1	9	468	A2M	N6-C6-N1	-8.31	100.15	118.35
1	9	623	A2M	C5-C6-N6	7.52	139.80	123.43

There are no chirality outliers.

5 of 43 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	9	28	A2M	C3'-C4'-C5'-O5'
1	9	38	OMC	C3'-C4'-C5'-O5'
1	9	38	OMC	O4'-C4'-C5'-O5'
1	9	246	OMG	O4'-C4'-C5'-O5'
1	9	246	OMG	C3'-C4'-C5'-O5'

There are no ring outliers.

9 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	9	468	A2M	3	0
1	9	392	OMG	2	0
1	9	140	OMC	1	0
1	9	615	OMU	2	0
1	9	1013	OMU	1	0
1	9	1781	4AC	2	0
1	9	978	A2M	1	0
1	9	1789	MA6	2	0
1	9	473	OMC	2	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 18 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

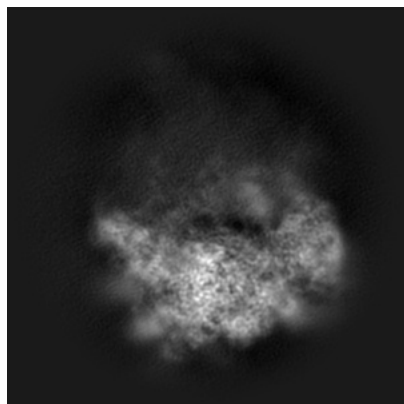
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-54789. These allow visual inspection of the internal detail of the map and identification of artifacts.

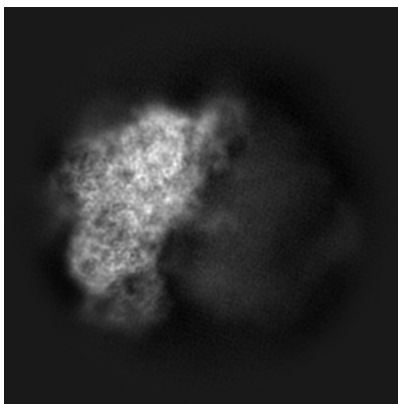
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

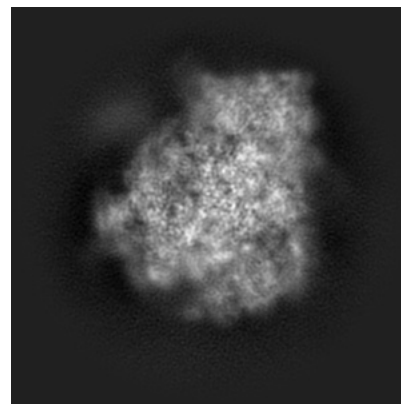
#### 6.1.1 Primary map



X

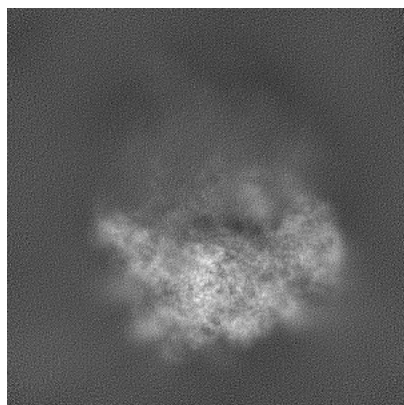


Y

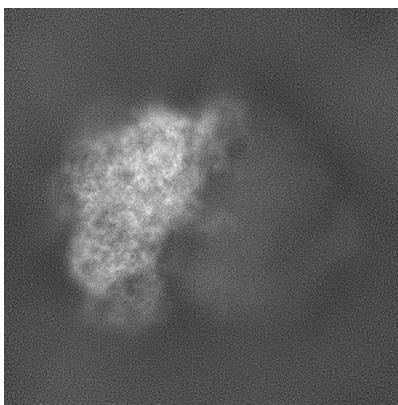


Z

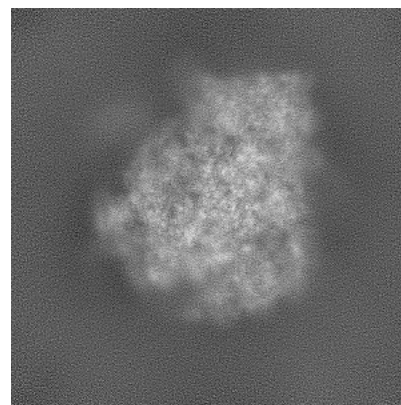
#### 6.1.2 Raw map



X



Y

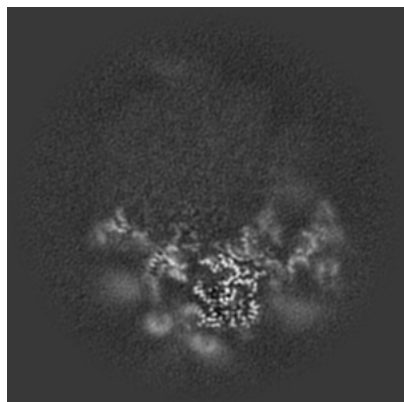


Z

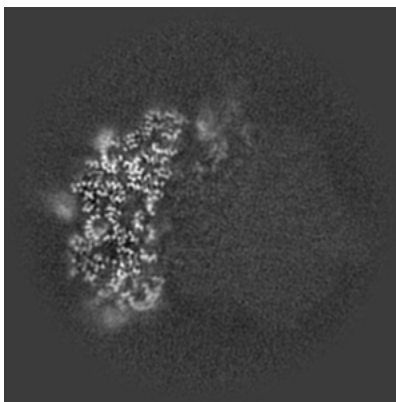
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

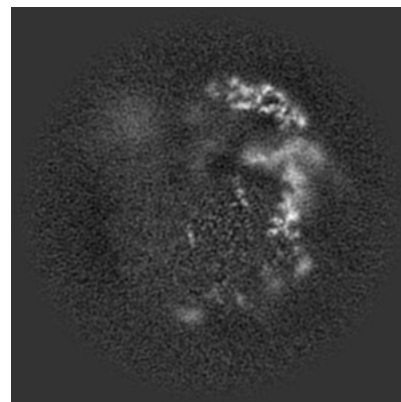
### 6.2.1 Primary map



X Index: 225

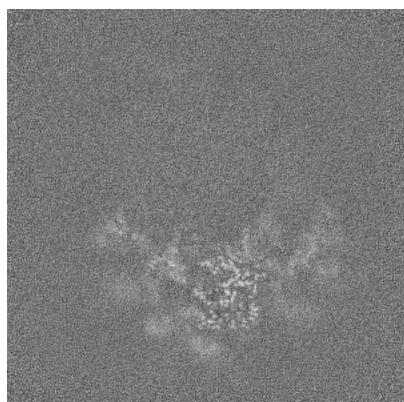


Y Index: 225

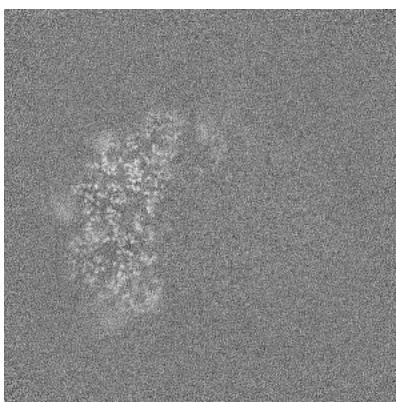


Z Index: 225

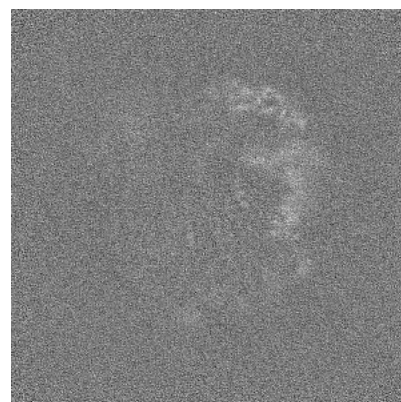
### 6.2.2 Raw map



X Index: 225



Y Index: 225

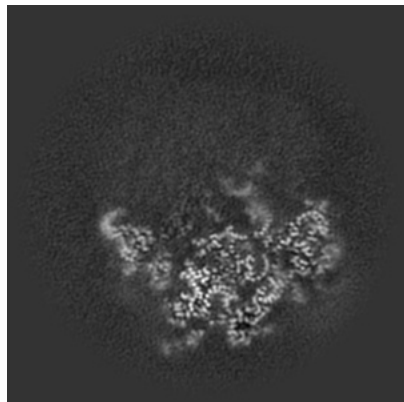


Z Index: 225

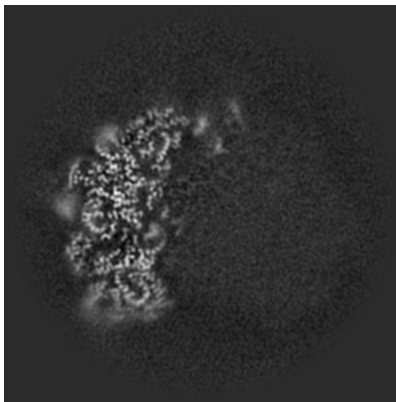
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

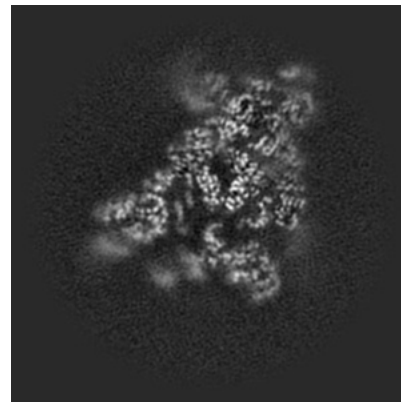
### 6.3.1 Primary map



X Index: 269

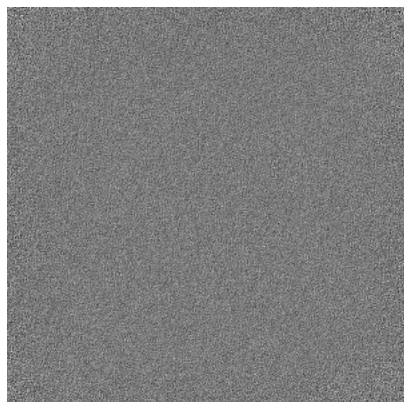


Y Index: 218

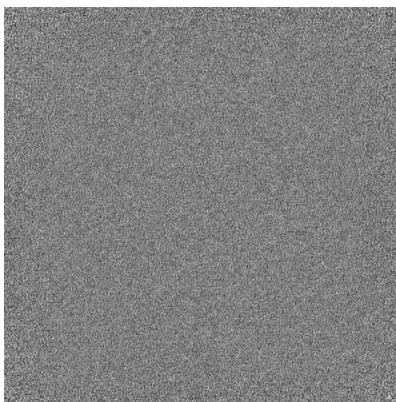


Z Index: 160

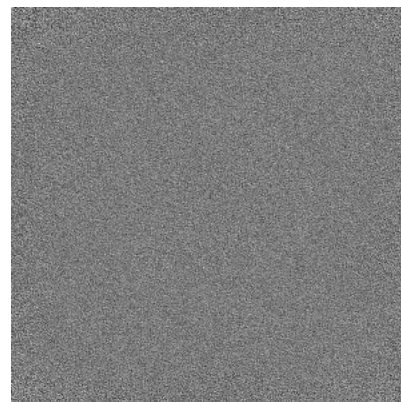
### 6.3.2 Raw map



X Index: 0



Y Index: 0

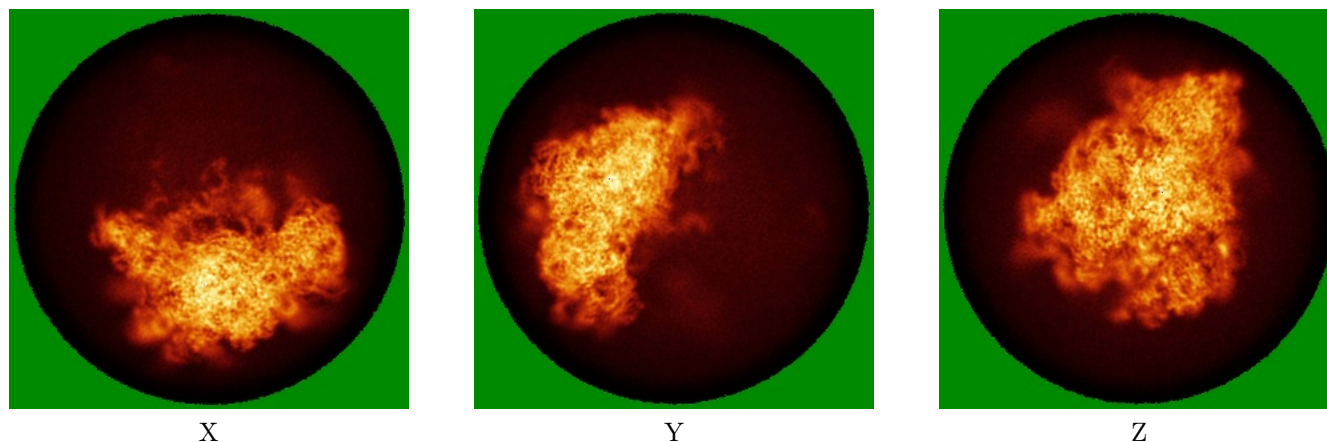


Z Index: 449

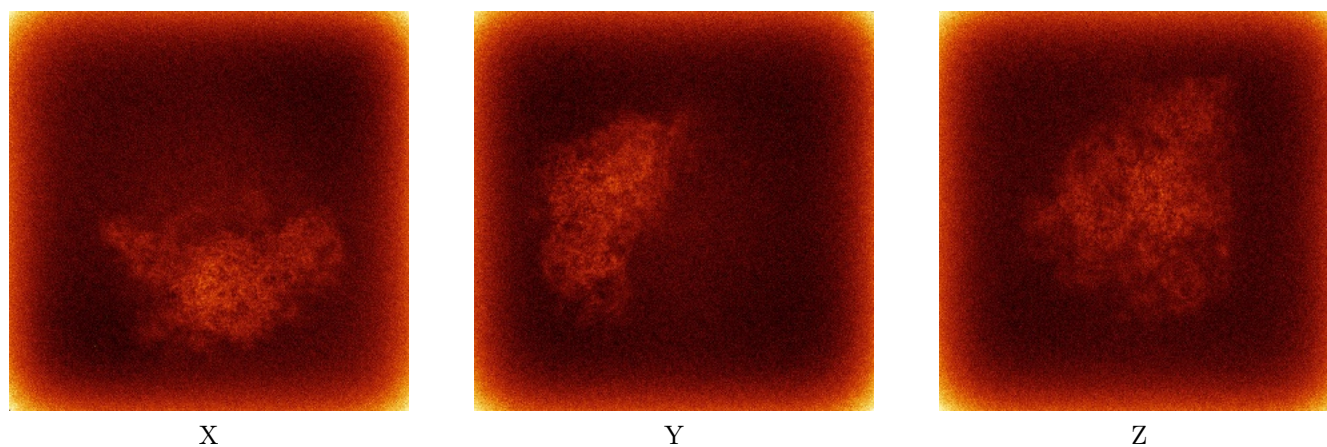
The images above show the largest variance slices of the map in three orthogonal directions.

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

### 6.4.1 Primary map



### 6.4.2 Raw map



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.

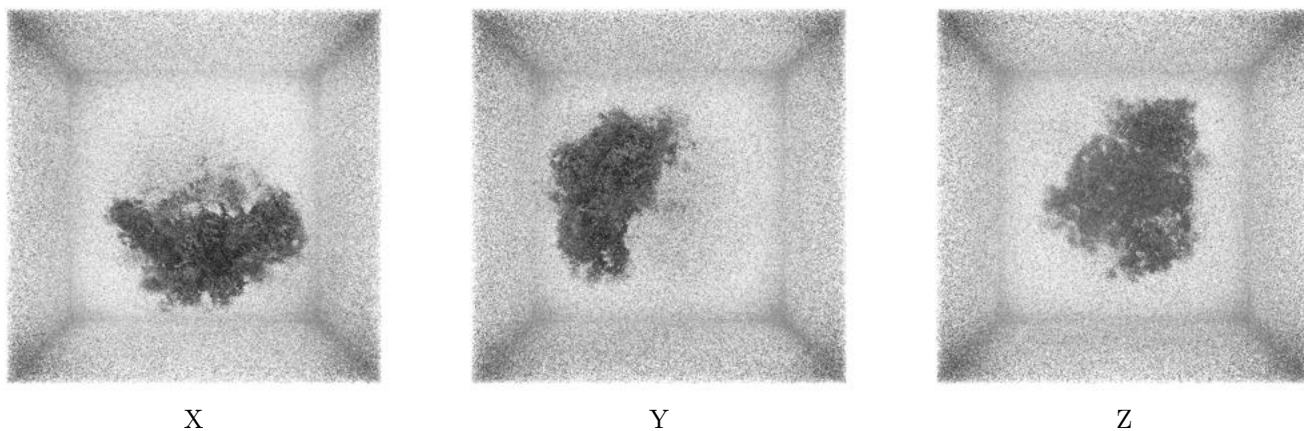
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



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

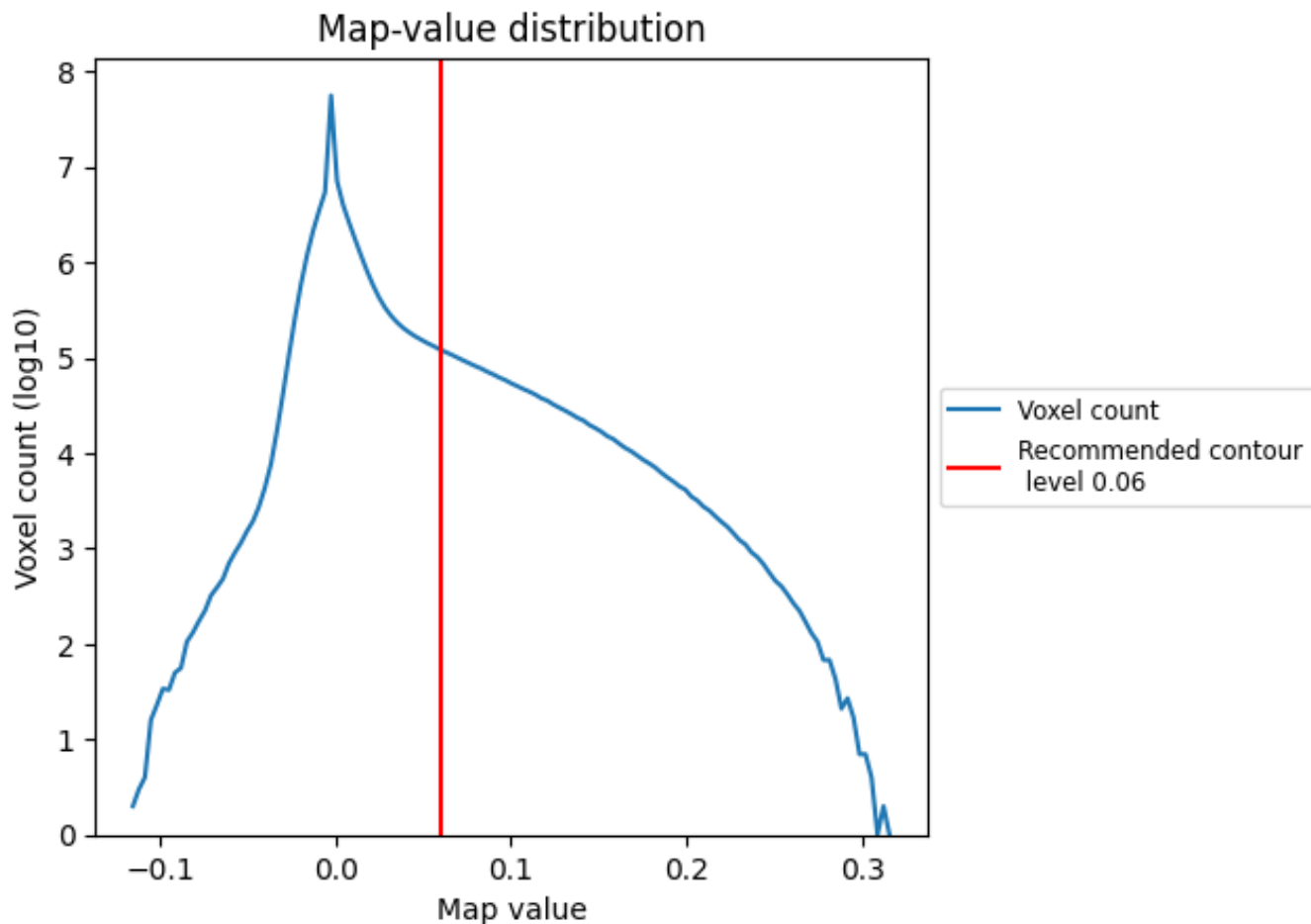
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

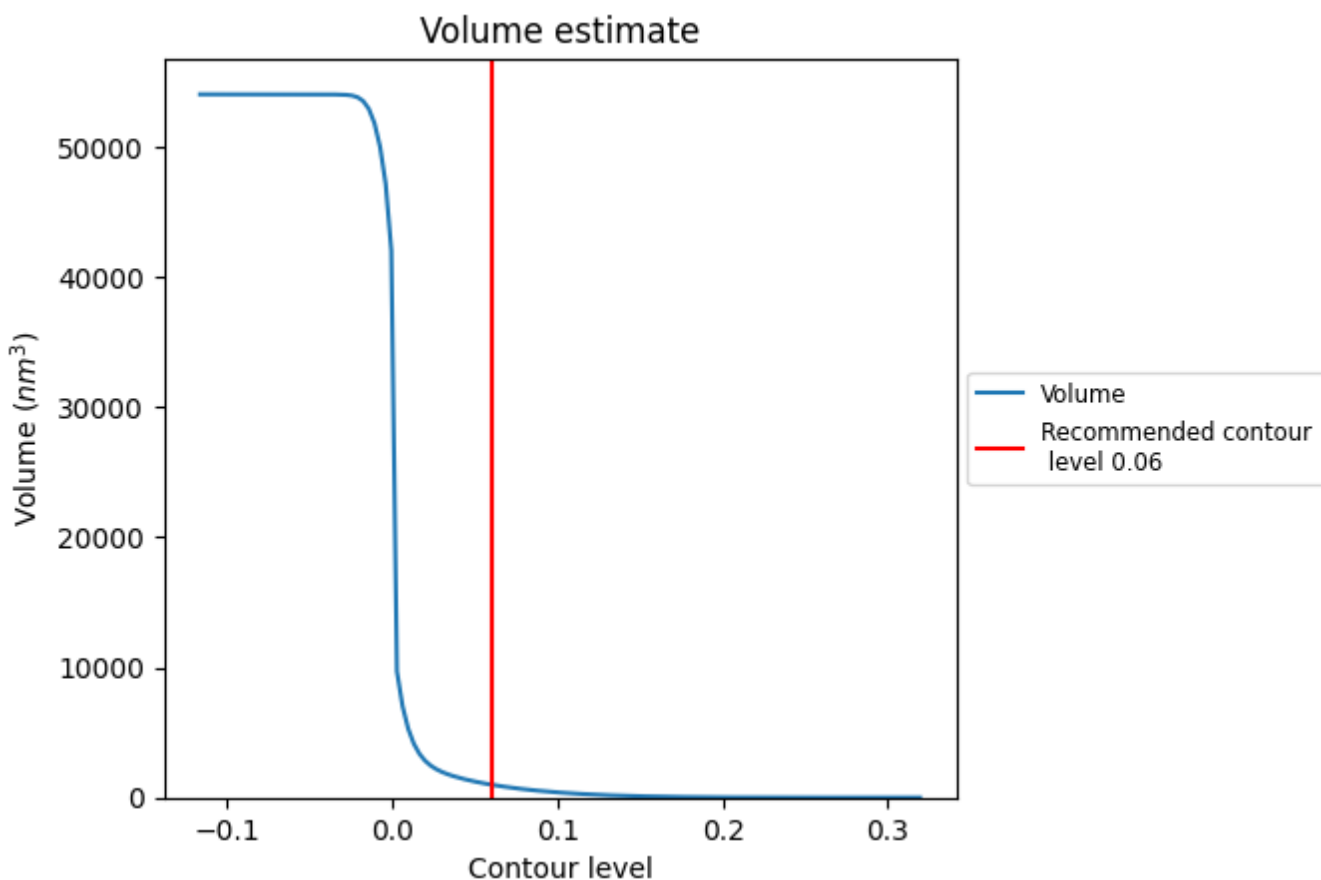
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

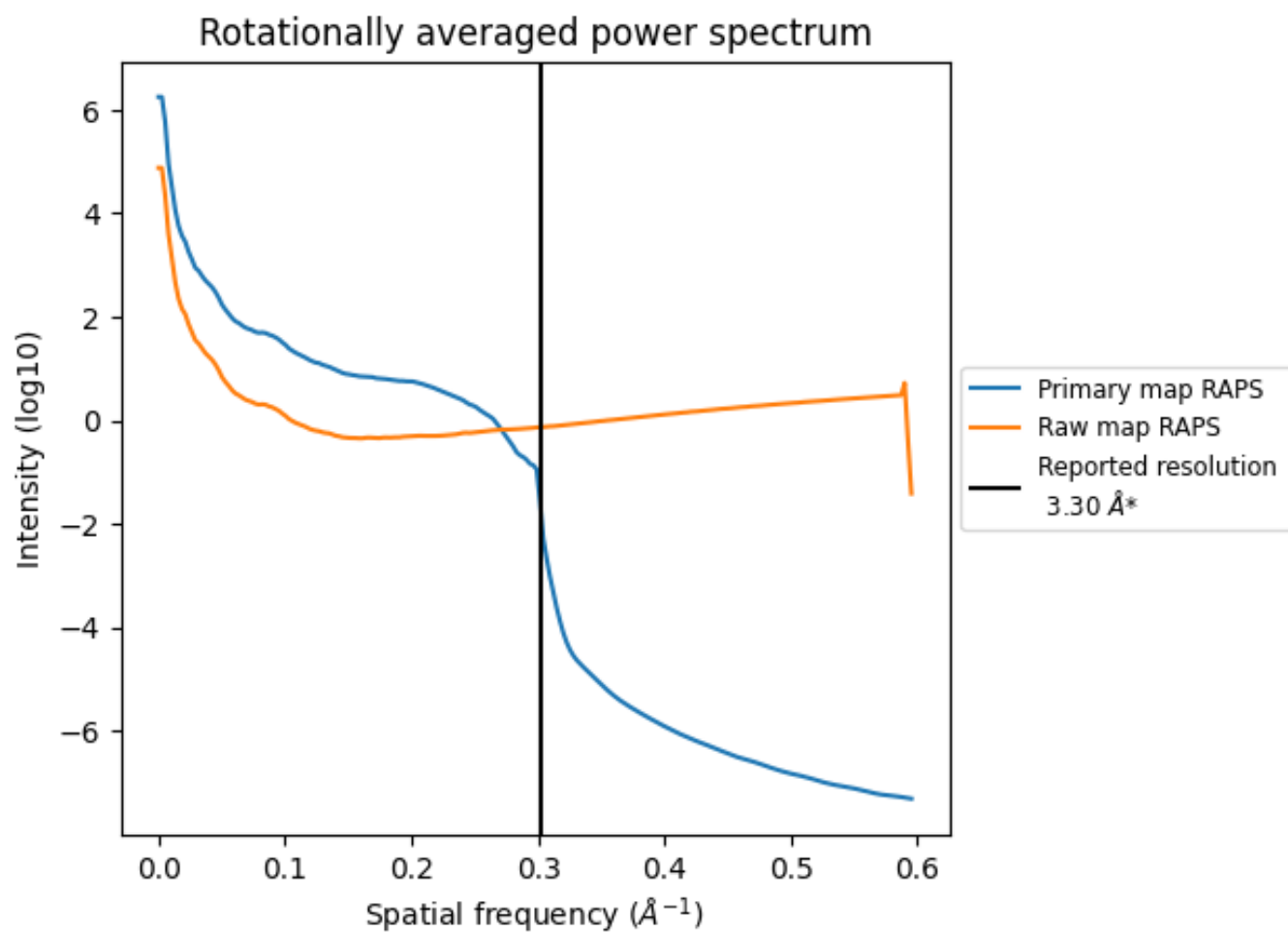
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 997 nm<sup>3</sup>; this corresponds to an approximate mass of 901 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 i

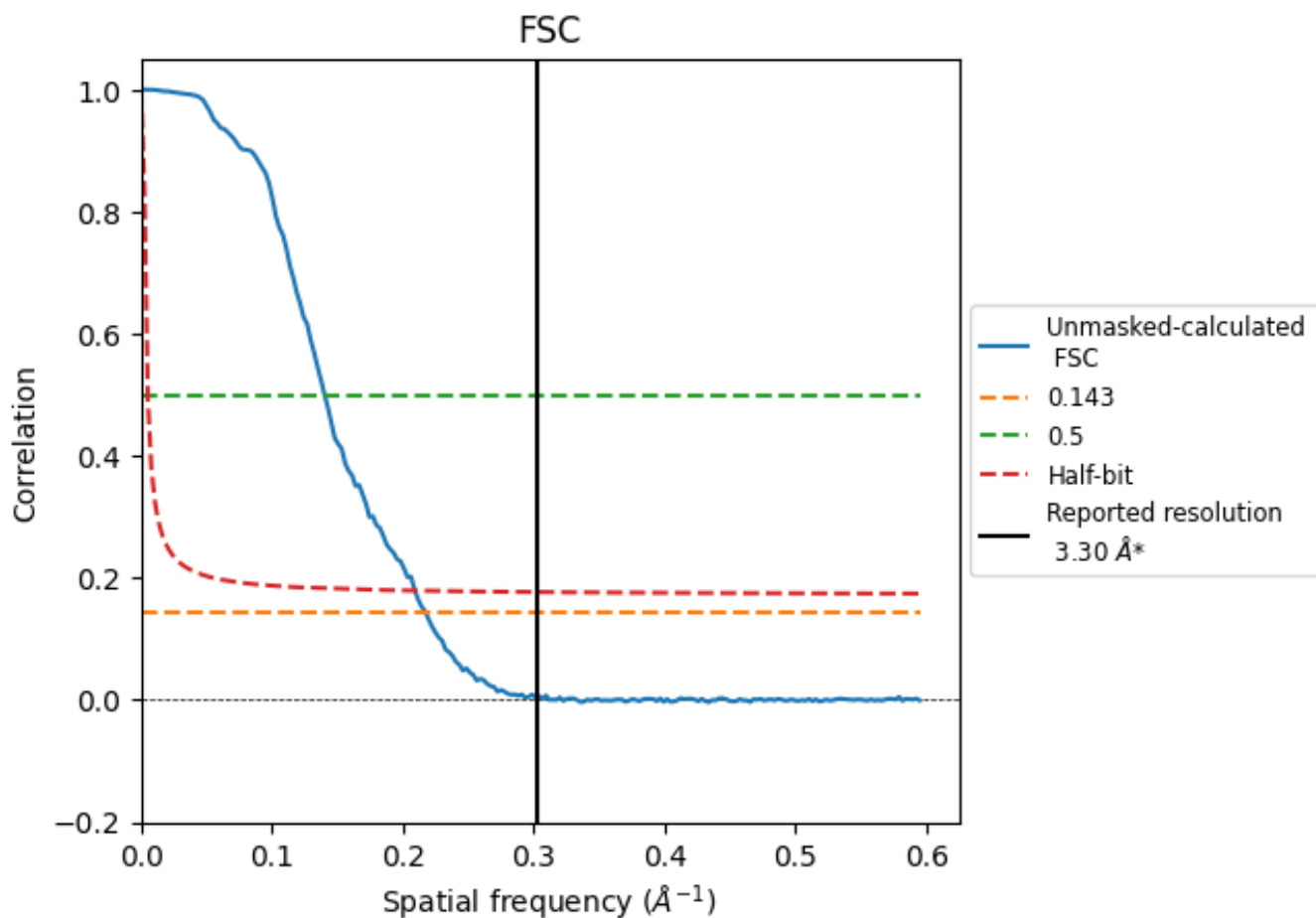


\*Reported resolution corresponds to spatial frequency of 0.303 Å<sup>-1</sup>

## 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 0.303 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

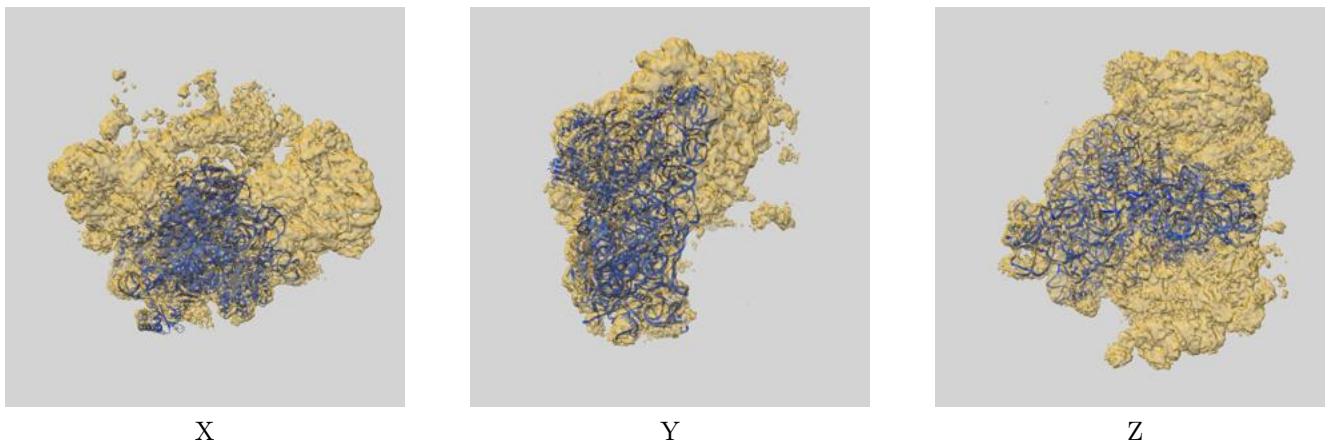
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.60	7.13	4.78

\*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 4.60 differs from the reported value 3.3 by more than 10 %

## 9 Map-model fit [i](#)

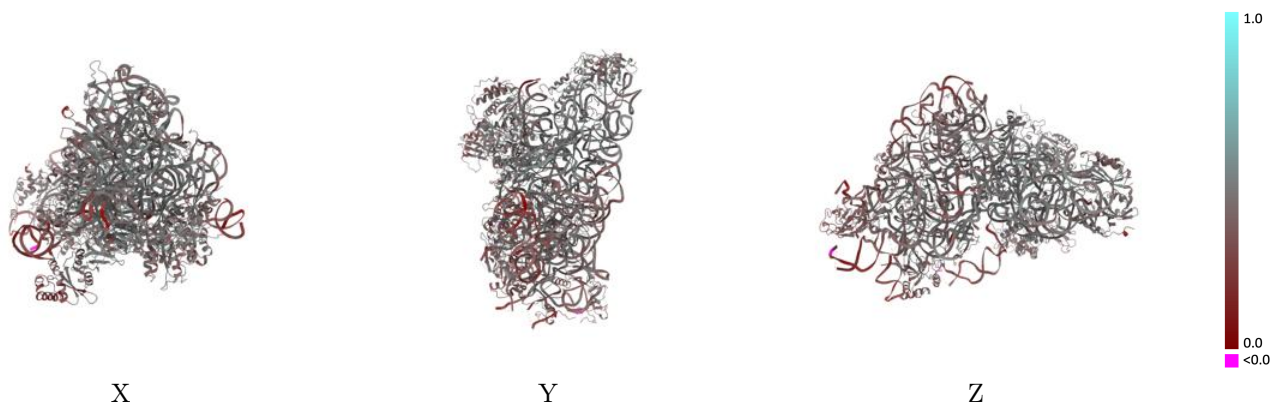
This section contains information regarding the fit between EMDB map EMD-54789 and PDB model 9SDP. Per-residue inclusion information can be found in section 3 on page 8.

### 9.1 Map-model overlay [i](#)



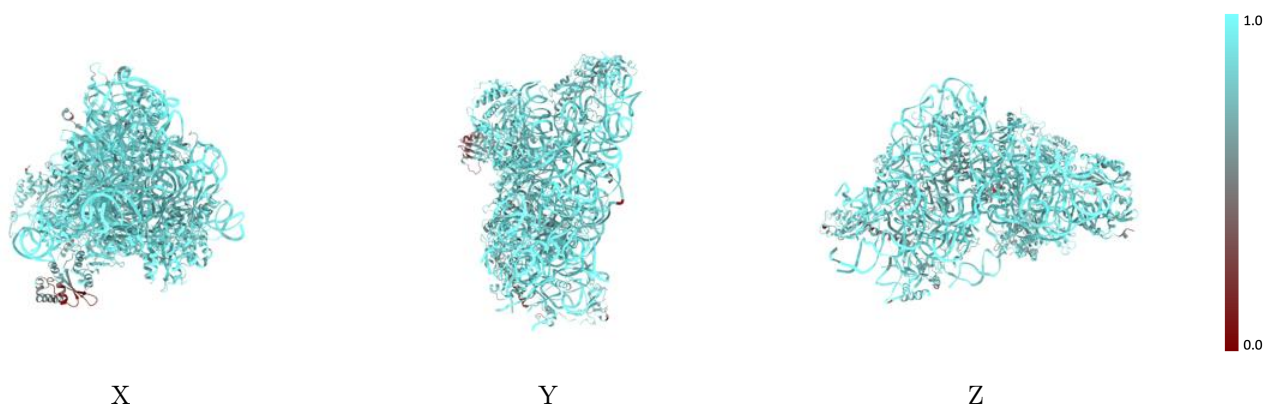
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



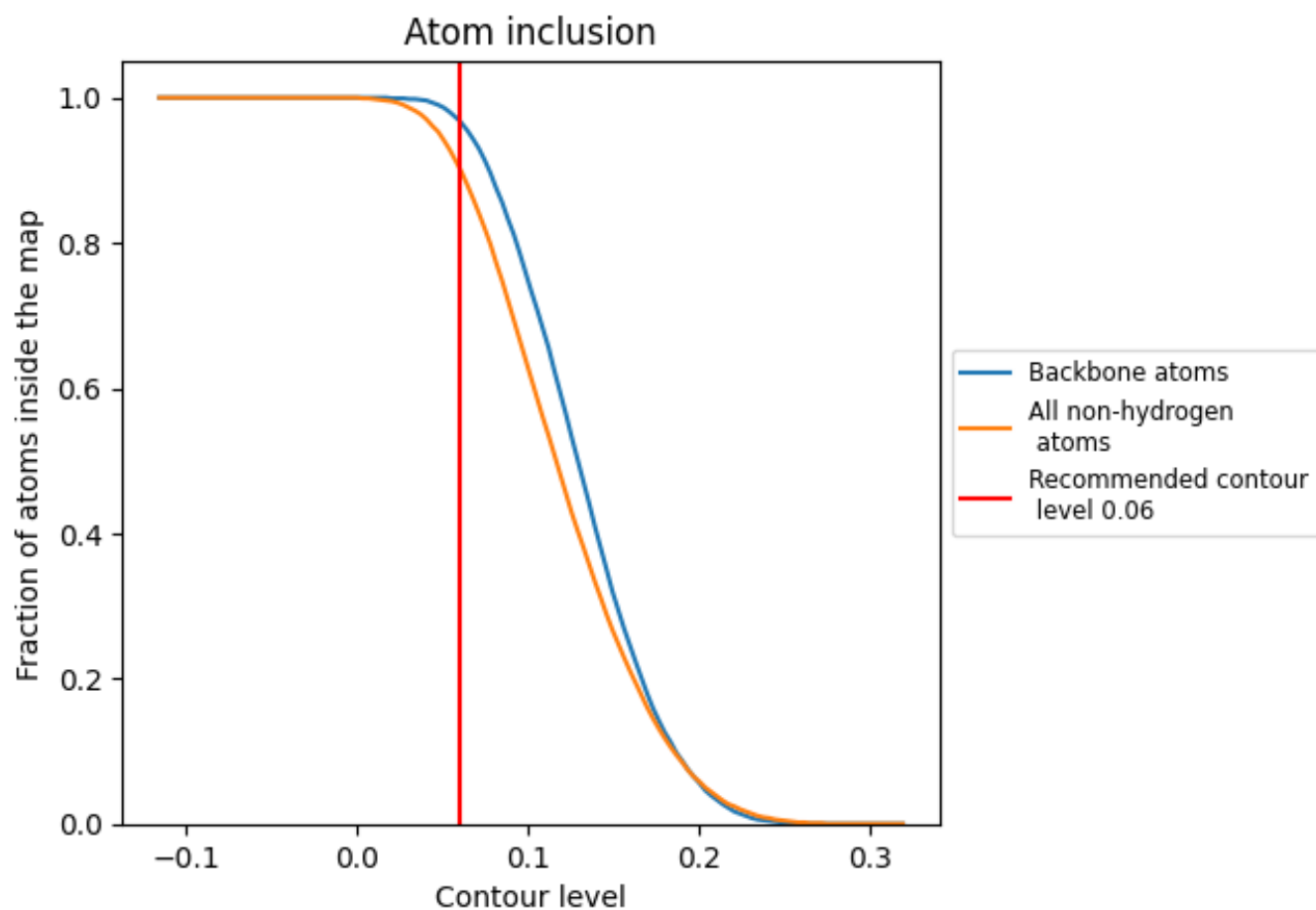
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).











































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 90% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9030	 0.4150
9	 0.9560	 0.4030
BA	 0.9010	 0.4310
BB	 0.8530	 0.4480
BC	 0.8990	 0.4600
BE	 0.7980	 0.4220
BG	 0.7220	 0.3700
BH	 0.6420	 0.3730
BI	 0.8720	 0.4160
BJ	 0.8570	 0.3990
BL	 0.9040	 0.4630
BN	 0.8990	 0.4550
BV	 0.9110	 0.4610
BW	 0.9160	 0.4850
BX	 0.8710	 0.4350
BY	 0.8620	 0.3670
BZ	 0.8440	 0.4720
Bb	 0.9140	 0.4410
Be	 0.8600	 0.3720
Cn	 0.7320	 0.4540
cS	 0.8780	 0.4540

