



wwPDB EM Validation Summary Report ⓘ

Apr 16, 2024 – 01:37 am BST

PDB ID : 7AFL
EMDB ID : EMD-11769
Title : Bacterial 30S ribosomal subunit assembly complex state D (multibody refinement for body domain of 30S ribosome)
Authors : Schedlbauer, A.; Iturrioz, I.; Ochoa-Lizarralde, B.; Diercks, T.; Kaminishi, T.; Capuni, R.; Astigarraga, E.; Gil-Carton, D.; Fucini, P.; Connell, S.
Deposited on : 2020-09-19
Resolution : 4.20 Å (reported)
Based on initial model : 4YBB

This is a wwPDB EM Validation Summary 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:

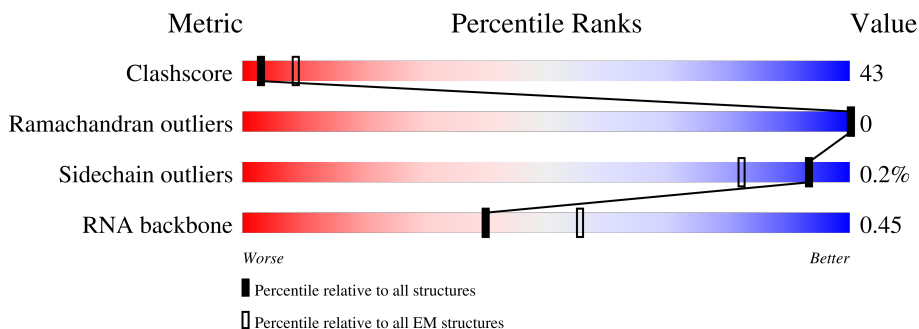
EMDB validation analysis : 0.0.1.dev92
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

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 4.20 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.



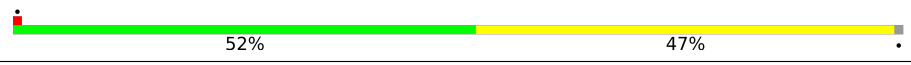

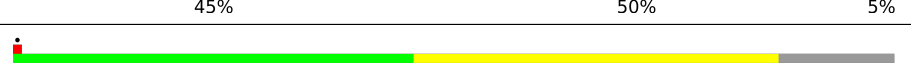
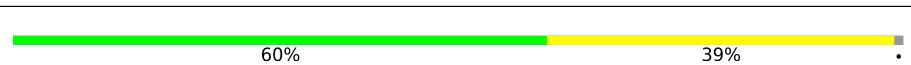
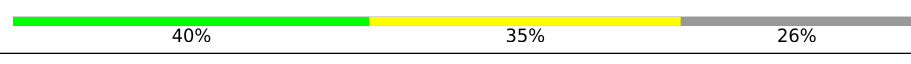


Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	1542	7% (red), 45% (yellow), 16% (orange), 32% (grey)
2	D	206	57% (green), 42% (yellow)
3	E	167	49% (green), 45% (yellow), 7% (grey)
4	F	135	34% (green), 44% (yellow), 21% (grey)
5	H	130	55% (green), 44% (yellow), . (grey)
6	K	129	52% (green), 39% (yellow), 9% (grey)
7	L	124	7% (red), 65% (green), 34% (yellow), . (grey)

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Mol	Chain	Length	Quality of chain
8	O	89	
9	P	82	
10	Q	84	
11	R	75	
12	T	87	
13	V	133	
14	X	151	

2 Entry composition [i](#)

There are 15 unique types of molecules in this entry. The entry contains 34322 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 16SrRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	A	1054	22654	10103	4173	7324	1054	0	0

- Molecule 2 is a protein called 30S ribosomal protein S4.

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

- Molecule 3 is a protein called 30S ribosomal protein S5.

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

- Molecule 4 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	F	106	862	545	156	154	7	0	0

- Molecule 5 is a protein called 30S ribosomal protein S8.

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

- Molecule 6 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	K	117	877	540	174	160	3	0	0

- Molecule 7 is a protein called 30S ribosomal protein S12.

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

- Molecule 8 is a protein called 30S ribosomal protein S15.

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

- Molecule 9 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	P	82	649	406	128	114	1	0	0

- Molecule 10 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	Q	80	648	411	121	113	3	0	0

- Molecule 11 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	R	65	535	339	100	95	1	0	0

- Molecule 12 is a protein called 30S ribosomal protein S20.

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

- Molecule 13 is a protein called Ribosome-binding factor A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	V	99	783	497	137	144	5	0	0

- Molecule 14 is a protein called Ribosome maturation factor RimP.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
14	X	151	1174	742	197	230	5	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	151	ALA	-	expression tag	UNP A0A0J3VRH1

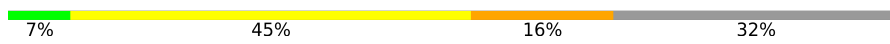
- Molecule 15 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

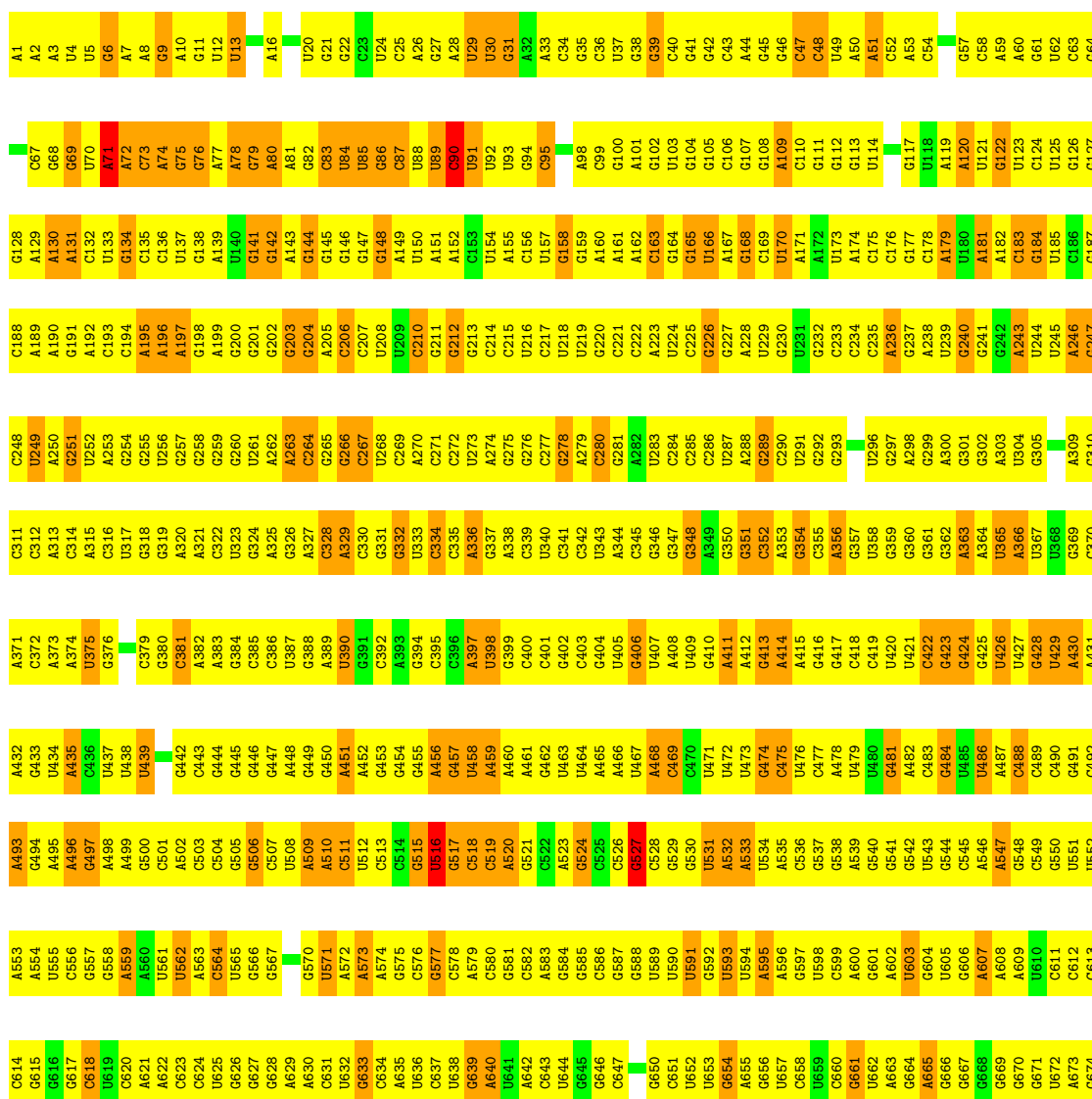
Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
15	A	25	25	25	0

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: 16SrRNA

Chain A: 



A675	A676	A677	U678	C679	A680	A681	A682	A683	A684	A685	A686	A687	A688	A689	A690	A691	A692	A693	A694	A695	A696	A697	A698	A699	A700	A701	A702	A703	A704	A705	A706	A707	A708	A709	A710	A711	A712	A713	A714	A715	A716	A717	A718	A719	A720	A721	A722	A723	A724	A725	A726	A727	A728	A729	A730	A731	A732	A733	A734																
C735	C736	C737	C738	C739	C740	C741	C742	C743	C744	C745	C746	C747	C748	C749	C750	C751	C752	C753	C754	C755	C756	C757	C758	C759	C760	C761	C762	C763	C764	C765	C766	C767	C768	C769	C770	C771	C772	C773	C774	C775	C776	C777	C778	C779	C780	C781	C782	C783	C784	C785	C786	C787	C788	C789	C790	C791	C792	C793	C794																
C795	C796	C797	C798	C799	C800	C801	C802	C803	C804	C805	C806	C807	C808	C809	C810	C811	C812	C813	C814	C815	C816	C817	C818	C819	C820	C821	C822	C823	C824	C825	C826	C827	C828	C829	C830	C831	C832	C833	C834	C835	C836	C837	C838	C839	C840	C841	C842	C843	C844	C845	C846	C847	C848	C849	C850	C851	C852	C853	C854																
U855	C856	C857	C858	C859	C860	C861	C862	C863	C864	C865	C866	C867	C868	C869	C870	C871	C872	C873	C874	C875	C876	C877	C878	C879	C880	C881	C882	C883	C884	C885	C886	C887	C888	C889	C890	C891	C892	C893	C894	C895	C896	C897	C898	C899	C900	C901	C902	C903	C904	C905	C906	C907	C908	C909	C910	C911	C912	C913	C914	C915	C916	C917	C918	C919	C920	C921	C922	C923	C924						
C925	C926	C927	C928	C929	C930	C931	C932	C933	C934	C935	C936	C937	C938	C939	C940	C941	C942	C943	C944	C945	C946	C947	C948	C949	C950	C951	C952	C953	C954	C955	C956	C957	C958	C959	C960	C961	C962	C963	C964	C965	C966	C967	C968	C969	C970	C971	C972	C973	C974	C975	C976	C977	C978	C979	C980	C981	C982	C983	C984	C985	C986	C987	C988	C989	C990	C991	C992	C993	C994	C995	C996	C997	C998	C999	C1000

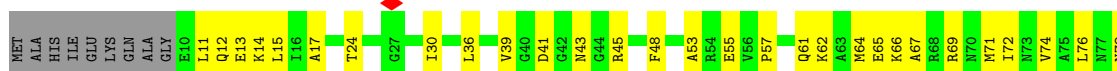
● Molecule 2: 30S ribosomal protein S4



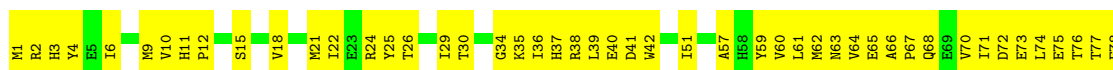
ME1	A2	R3	K8	L11	S12	R13	R14	E15	G16	T17	D18	L19	F20	H20	L21	K22	I28	K31	C32	K33	I34	E35	P38	G39	Q40	R47	L48	S49	D50	Y51	G52	V53	Q54	L55	R56	E57	K58	Q59	R62	R63	E69	R70	K83	L91	E95	D99	N100	Y101	
V102	F103	R104	M105	G106	F107	T110	E113	A114	R115	O116	L117	V118	S119	H120	L121	K121	A122	I123	M124	V125	R128	V129	I132	V137	S138	D141	V142	V143	S144	Q152	S153	R154	V155	A158	L159	E160	P168	T169	W170	L171	F172	V173	D174	K177	M178	E179	G180	T181	F182



- Molecule 3: 30S ribosomal protein S5



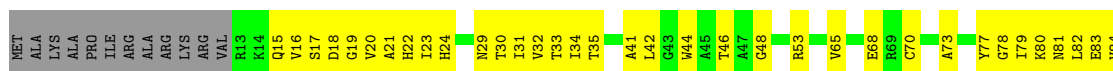
- Molecule 4: 30S ribosomal protein S6



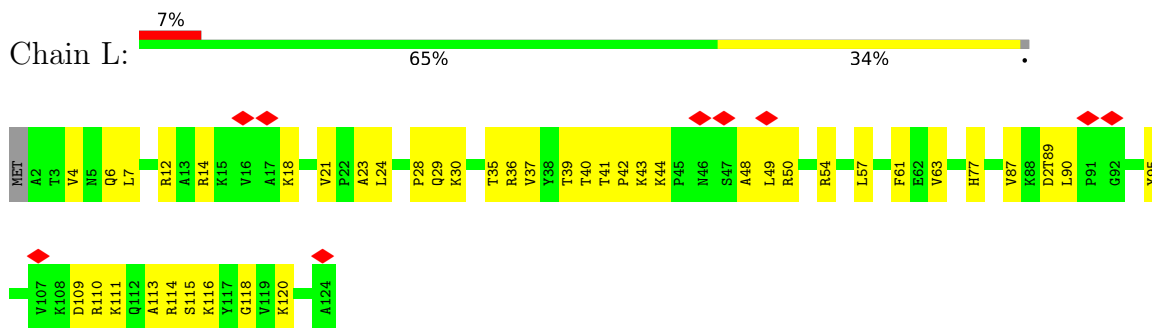
- Molecule 5: 30S ribosomal protein S8



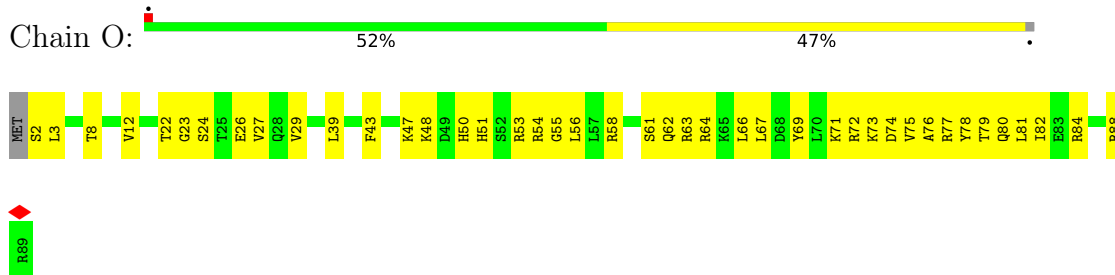
- Molecule 6: 30S ribosomal protein S11



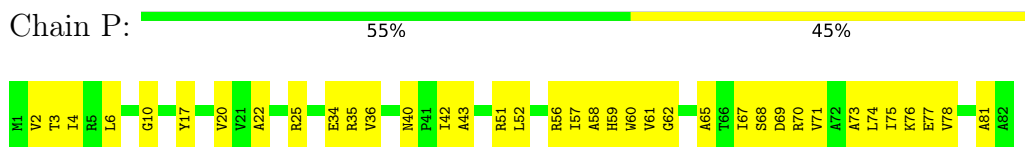
- Molecule 7: 30S ribosomal protein S12



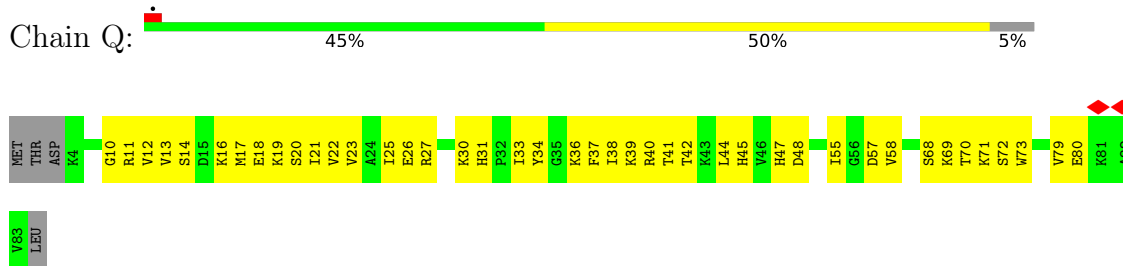
• Molecule 8: 30S ribosomal protein S15



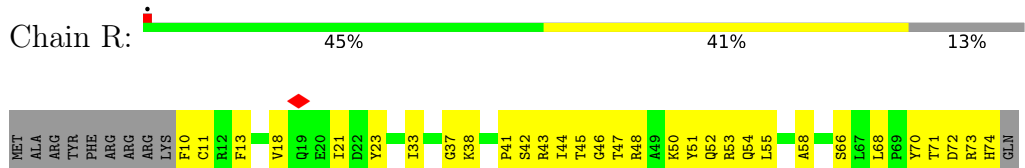
• Molecule 9: 30S ribosomal protein S16



• Molecule 10: 30S ribosomal protein S17



• Molecule 11: 30S ribosomal protein S18

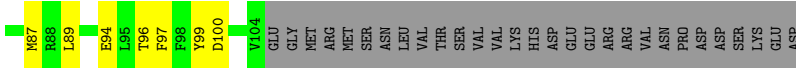
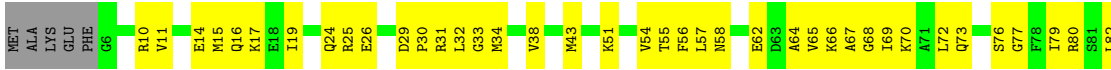


• Molecule 12: 30S ribosomal protein S20

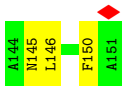
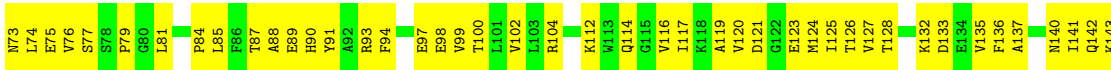
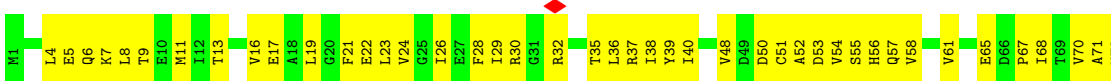




• Molecule 13: Ribosome-binding factor A



• Molecule 14: Ribosome maturation factor RimP



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	4573	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	46.1	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.098	Depositor
Minimum map value	-0.029	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.018	Depositor
Map size (Å)	416.64, 416.64, 416.64	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.085, 1.085, 1.085	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: G7M, MG, 2MG, D2T, PSU, MA6

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	1.03	1/25238 (0.0%)	1.04	9/39369 (0.0%)
2	D	0.56	0/1665	0.62	0/2227
3	E	0.57	0/1165	0.64	0/1568
4	F	0.51	0/881	0.63	0/1189
5	H	0.58	0/989	0.67	0/1326
6	K	0.58	0/893	0.70	0/1205
7	L	0.58	0/960	0.66	0/1286
8	O	0.52	0/722	0.59	0/964
9	P	0.61	0/659	0.64	0/884
10	Q	0.59	0/657	0.65	0/881
11	R	0.56	0/544	0.64	0/731
12	T	0.54	0/676	0.61	0/895
13	V	0.57	0/792	0.68	1/1062 (0.1%)
14	X	0.50	0/1192	0.62	0/1619
All	All	0.91	1/37033 (0.0%)	0.94	10/55206 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	571	U	O3'-P	-11.48	1.47	1.61

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1431	A	C8-N9-C4	5.78	108.11	105.80
13	V	30	PRO	CA-N-CD	-5.78	103.40	111.50
1	A	90	C	O4'-C1'-N1	5.44	112.55	108.20
1	A	71	A	C8-N9-C4	5.35	107.94	105.80
1	A	439	U	N3-C2-O2	-5.33	118.47	122.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	22654	0	11396	1524	0
2	D	1643	0	1707	89	0
3	E	1152	0	1196	76	0
4	F	862	0	864	62	0
5	H	979	0	1031	71	0
6	K	877	0	887	64	0
7	L	957	0	1017	59	0
8	O	714	0	734	44	0
9	P	649	0	666	35	0
10	Q	648	0	691	44	0
11	R	535	0	552	36	0
12	T	670	0	719	34	0
13	V	783	0	816	70	0
14	X	1174	0	1174	95	0
15	A	25	0	0	0	0
All	All	34322	0	23450	2139	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 43.

The worst 5 of 2139 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)
13:V:32:LEU:CD1	13:V:34:MET:H	1.67	1.08
7:L:37:VAL:HG22	14:X:48:VAL:CG2	1.84	1.07
13:V:32:LEU:HD12	13:V:33:GLY:N	1.78	0.98
13:V:32:LEU:HD12	13:V:34:MET:H	1.24	0.96
13:V:11:VAL:HG23	13:V:43:MET:CE	1.95	0.95

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	D	203/206 (98%)	191 (94%)	12 (6%)	0	100	100
3	E	154/167 (92%)	145 (94%)	9 (6%)	0	100	100
4	F	104/135 (77%)	97 (93%)	7 (7%)	0	100	100
5	H	127/130 (98%)	118 (93%)	9 (7%)	0	100	100
6	K	115/129 (89%)	107 (93%)	8 (7%)	0	100	100
7	L	120/124 (97%)	114 (95%)	6 (5%)	0	100	100
8	O	86/89 (97%)	82 (95%)	4 (5%)	0	100	100
9	P	80/82 (98%)	76 (95%)	4 (5%)	0	100	100
10	Q	78/84 (93%)	72 (92%)	6 (8%)	0	100	100
11	R	63/75 (84%)	60 (95%)	3 (5%)	0	100	100
12	T	84/87 (97%)	78 (93%)	6 (7%)	0	100	100
13	V	97/133 (73%)	93 (96%)	4 (4%)	0	100	100
14	X	149/151 (99%)	140 (94%)	9 (6%)	0	100	100
All	All	1460/1592 (92%)	1373 (94%)	87 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	D	172/173 (99%)	171 (99%)	1 (1%)	86	92

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	E	119/126 (94%)	118 (99%)	1 (1%)	81	89
4	F	92/116 (79%)	92 (100%)	0	100	100
5	H	104/105 (99%)	104 (100%)	0	100	100
6	K	90/99 (91%)	90 (100%)	0	100	100
7	L	102/103 (99%)	102 (100%)	0	100	100
8	O	76/77 (99%)	76 (100%)	0	100	100
9	P	65/65 (100%)	65 (100%)	0	100	100
10	Q	74/78 (95%)	74 (100%)	0	100	100
11	R	56/65 (86%)	56 (100%)	0	100	100
12	T	65/66 (98%)	65 (100%)	0	100	100
13	V	84/116 (72%)	84 (100%)	0	100	100
14	X	130/130 (100%)	130 (100%)	0	100	100
All	All	1229/1319 (93%)	1227 (100%)	2 (0%)	93	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	83	LYS
3	E	78	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
9	P	29	ASN
12	T	20	HIS
11	R	52	GLN
12	T	84	ASN
6	K	24	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1048/1542 (67%)	282 (26%)	10 (0%)

5 of 282 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	6	G
1	A	9	G
1	A	13	U
1	A	16	A
1	A	22	G

5 of 10 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	776	G
1	A	789	U
1	A	884	U
1	A	428	G
1	A	429	U

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

6 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	MA6	A	1518	1	18,26,27	1.02	1 (5%)	19,38,41	1.73	5 (26%)
1	2MG	A	1516	1	18,26,27	0.88	1 (5%)	16,38,41	1.45	4 (25%)
7	D2T	L	89	7	7,9,10	1.04	1 (14%)	6,11,13	1.74	1 (16%)
1	PSU	A	516	1	18,21,22	1.41	3 (16%)	22,30,33	2.22	6 (27%)
1	G7M	A	527	1	20,26,27	0.92	1 (5%)	17,39,42	1.10	1 (5%)
1	MA6	A	1519	1	18,26,27	0.99	1 (5%)	19,38,41	1.94	6 (31%)

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	MA6	A	1518	1	-	4/7/29/30	0/3/3/3
1	2MG	A	1516	1	-	2/5/27/28	0/3/3/3
7	D2T	L	89	7	-	3/7/12/14	-
1	PSU	A	516	1	-	2/7/25/26	0/2/2/2
1	G7M	A	527	1	-	1/3/25/26	0/3/3/3
1	MA6	A	1519	1	-	4/7/29/30	0/3/3/3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	516	PSU	C4-N3	-3.22	1.32	1.38
1	A	527	G7M	C8-N9	2.93	1.38	1.33
7	L	89	D2T	O-C	2.68	1.30	1.19
1	A	1519	MA6	C5-C4	2.54	1.47	1.40
1	A	1518	MA6	C5-C4	2.52	1.47	1.40

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	516	PSU	N1-C2-N3	6.61	122.62	115.13
1	A	516	PSU	C4-N3-C2	-4.61	119.70	126.34
1	A	1519	MA6	N1-C6-N6	4.12	121.40	117.06
1	A	1516	2MG	CM2-N2-C2	-3.72	115.64	123.86
1	A	516	PSU	O2-C2-N1	-3.65	118.77	122.79

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1516	2MG	N1-C2-N2-CM2
1	A	1516	2MG	N3-C2-N2-CM2
1	A	1518	MA6	O4'-C4'-C5'-O5'
1	A	1519	MA6	O4'-C4'-C5'-O5'
1	A	1519	MA6	C3'-C4'-C5'-O5'

There are no ring outliers.

5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1518	MA6	7	0
1	A	1516	2MG	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	516	PSU	5	0
1	A	527	G7M	1	0
1	A	1519	MA6	6	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 25 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

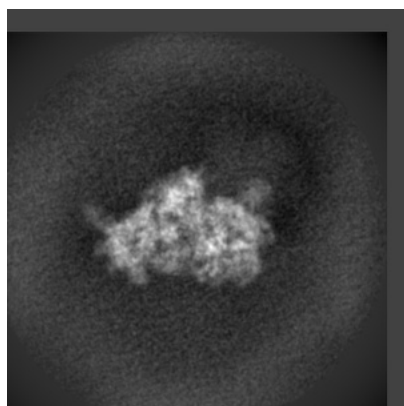
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11769. These allow visual inspection of the internal detail of the map and identification of artifacts.

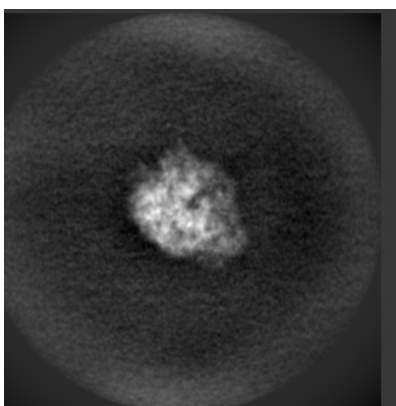
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

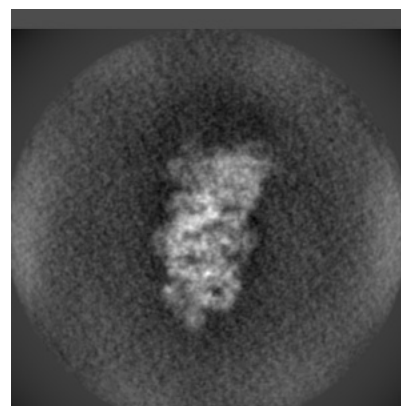
6.1.1 Primary map



X

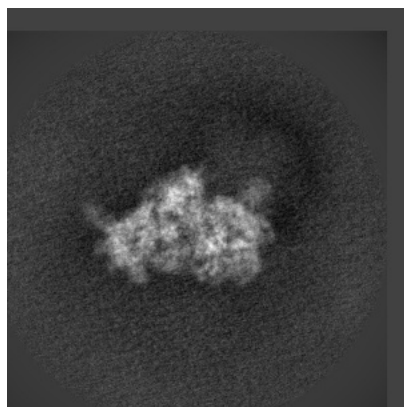


Y

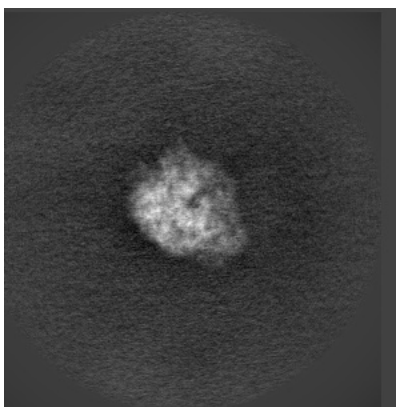


Z

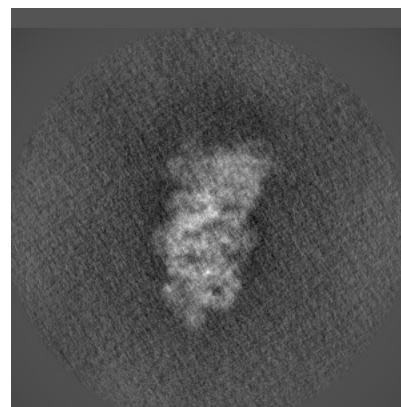
6.1.2 Raw map



X



Y

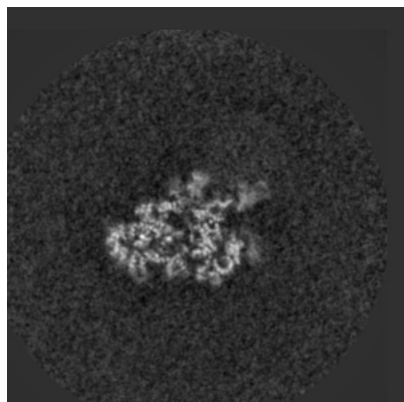


Z

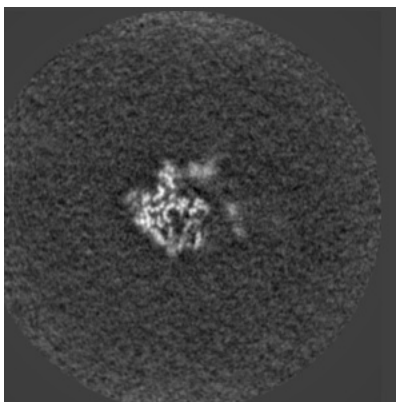
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

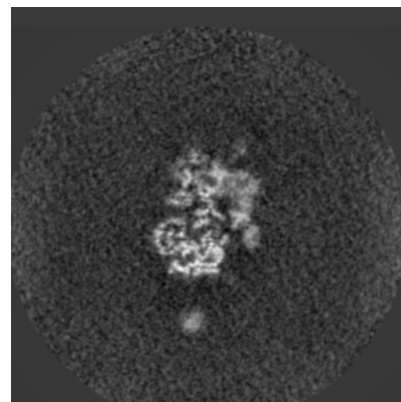
6.2.1 Primary map



X Index: 192

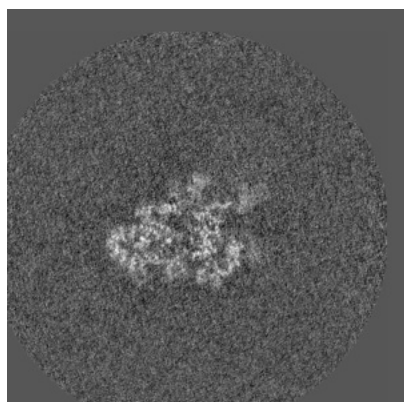


Y Index: 192

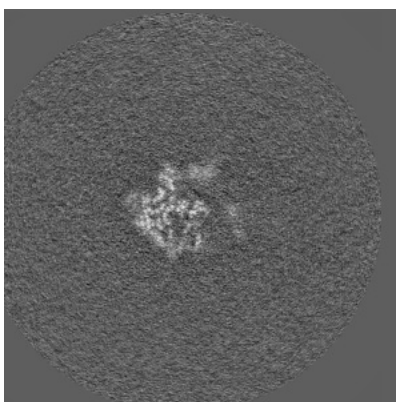


Z Index: 192

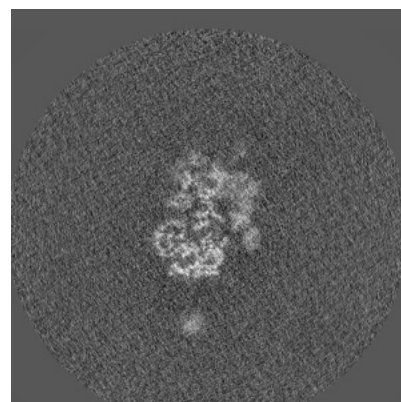
6.2.2 Raw map



X Index: 192



Y Index: 192

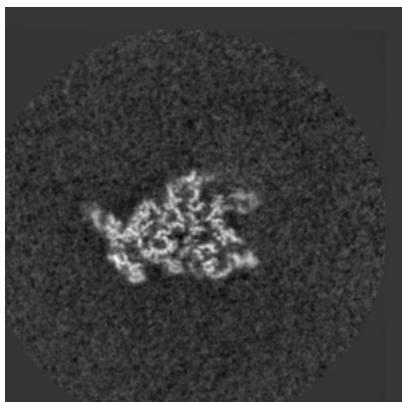


Z Index: 192

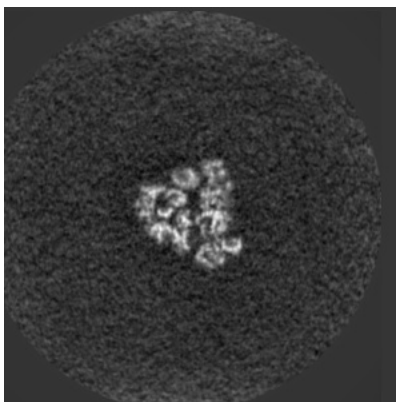
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

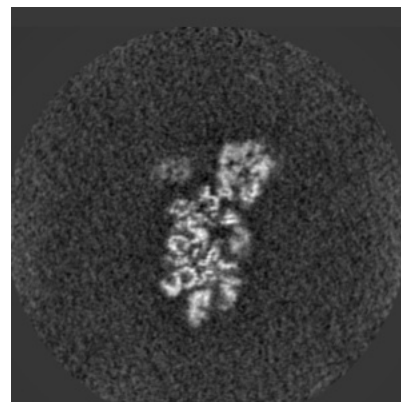
6.3.1 Primary map



X Index: 184

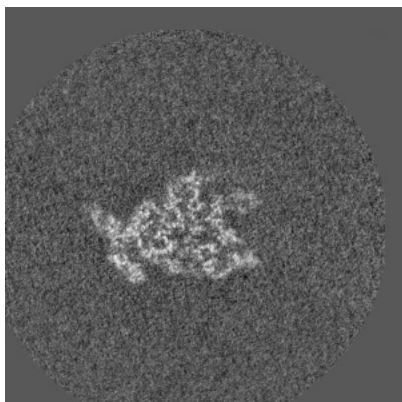


Y Index: 159

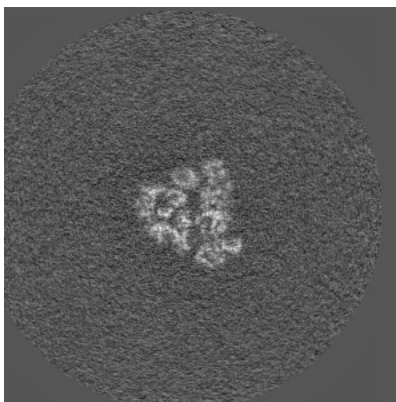


Z Index: 175

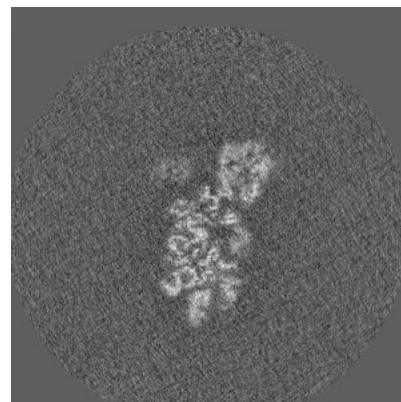
6.3.2 Raw map



X Index: 184



Y Index: 159

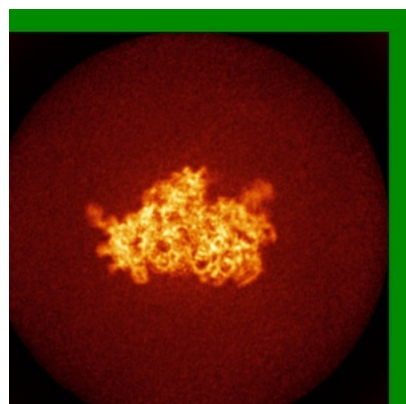


Z Index: 175

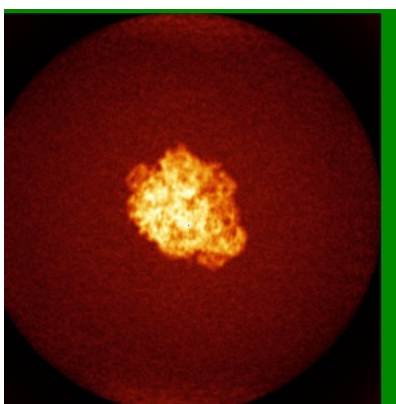
The images above show the largest variance slices of the map in three orthogonal directions.

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

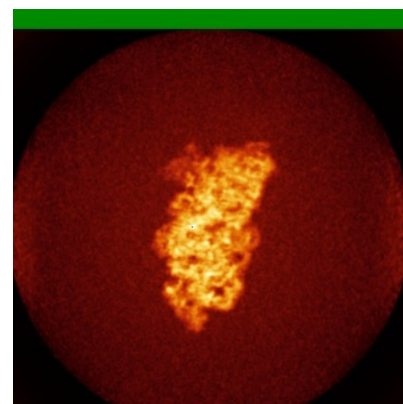
6.4.1 Primary map



X

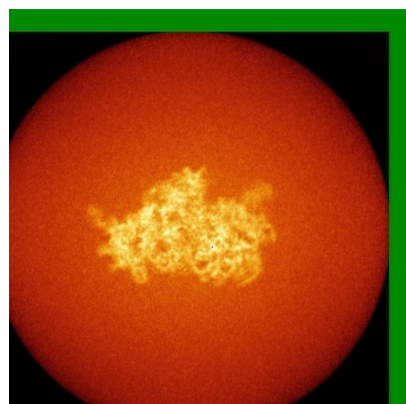


Y

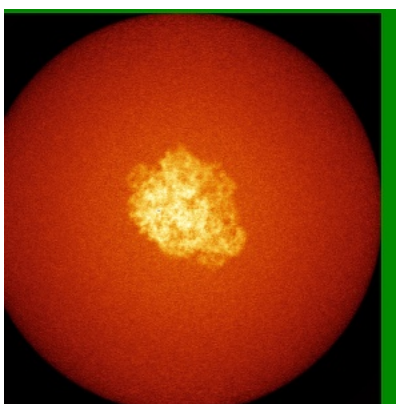


Z

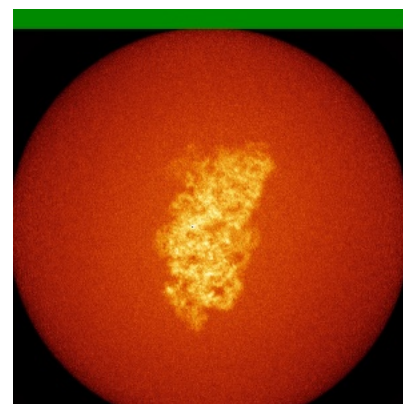
6.4.2 Raw map



X



Y

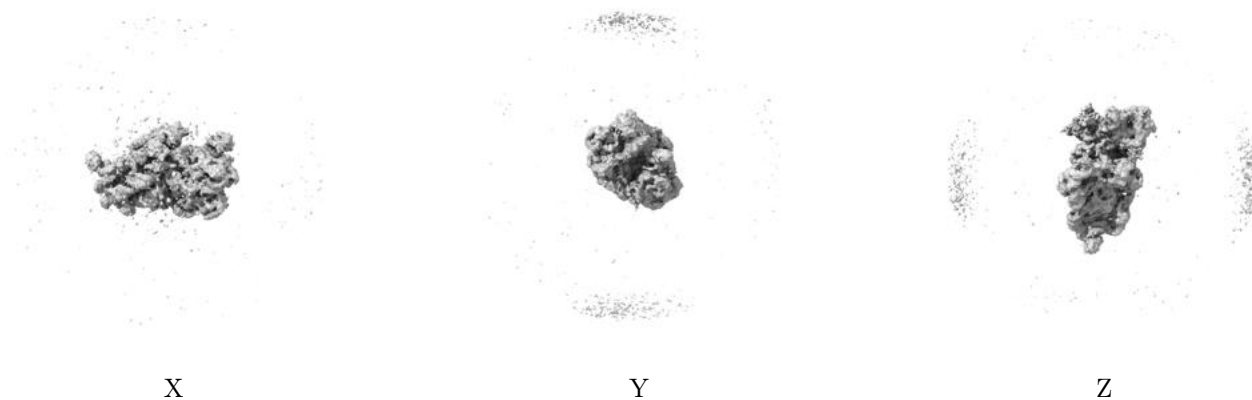


Z

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.

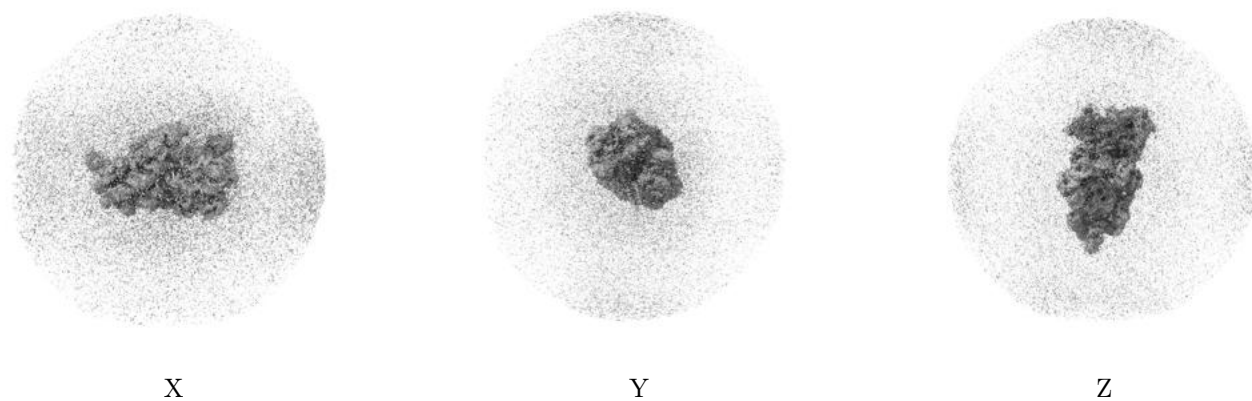
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.018. 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.

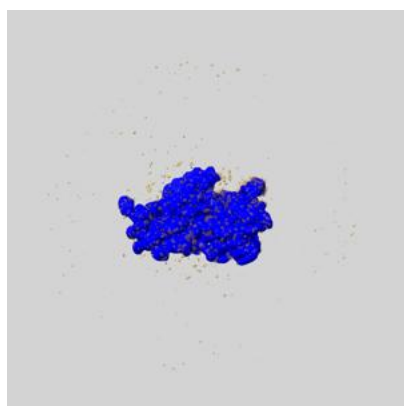
6.6 Mask visualisation [i](#)

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

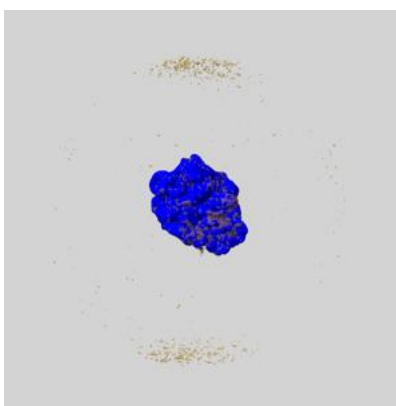
A mask typically either:

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

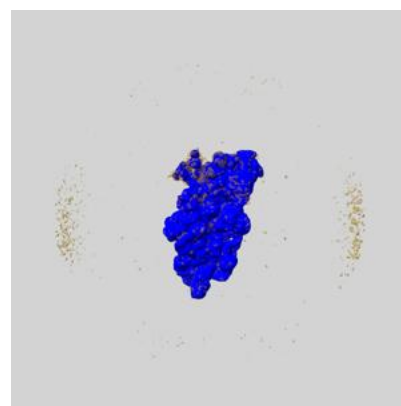
6.6.1 emd_11769_msk_1.map [i](#)



X



Y

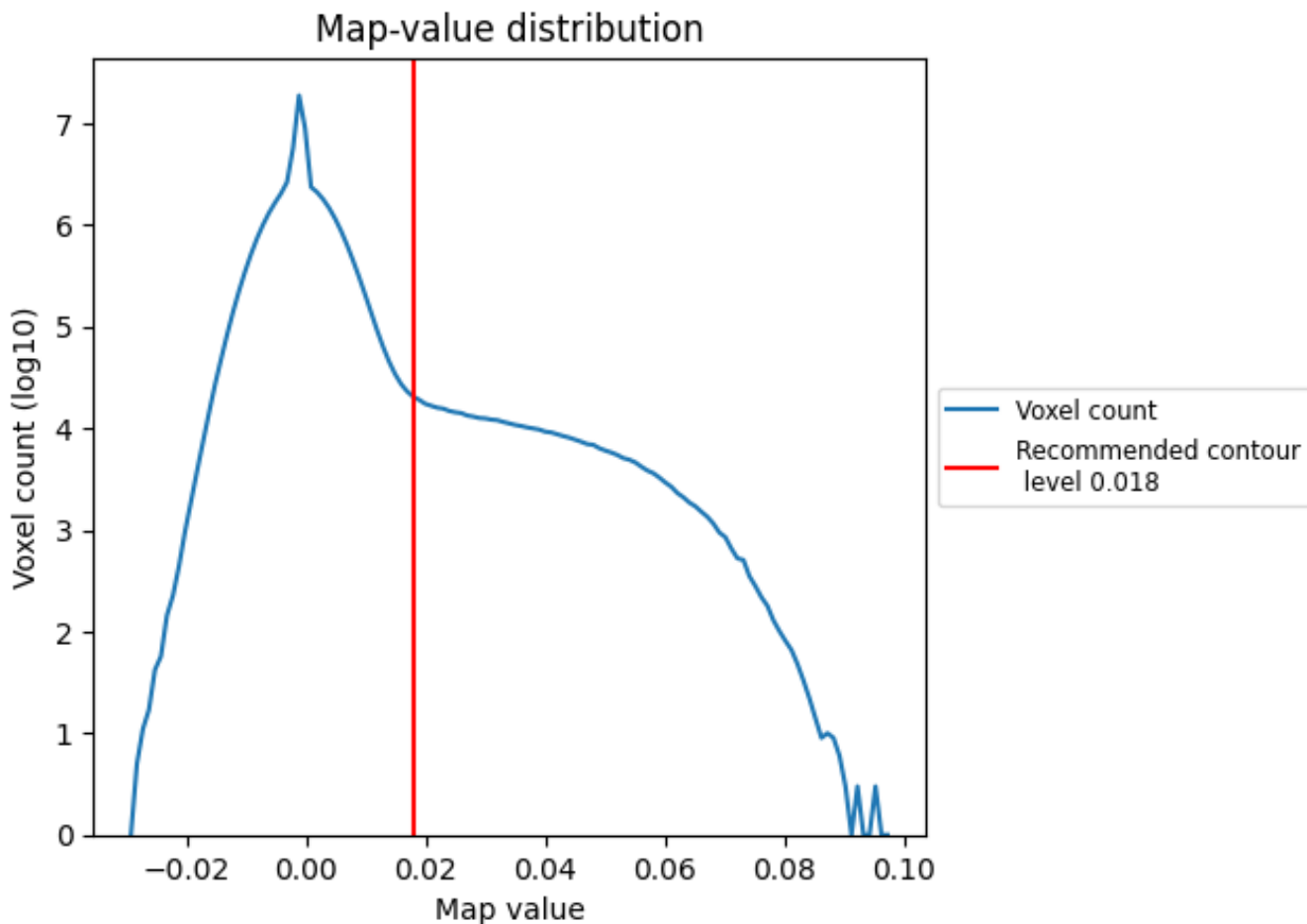


Z

7 Map analysis [i](#)

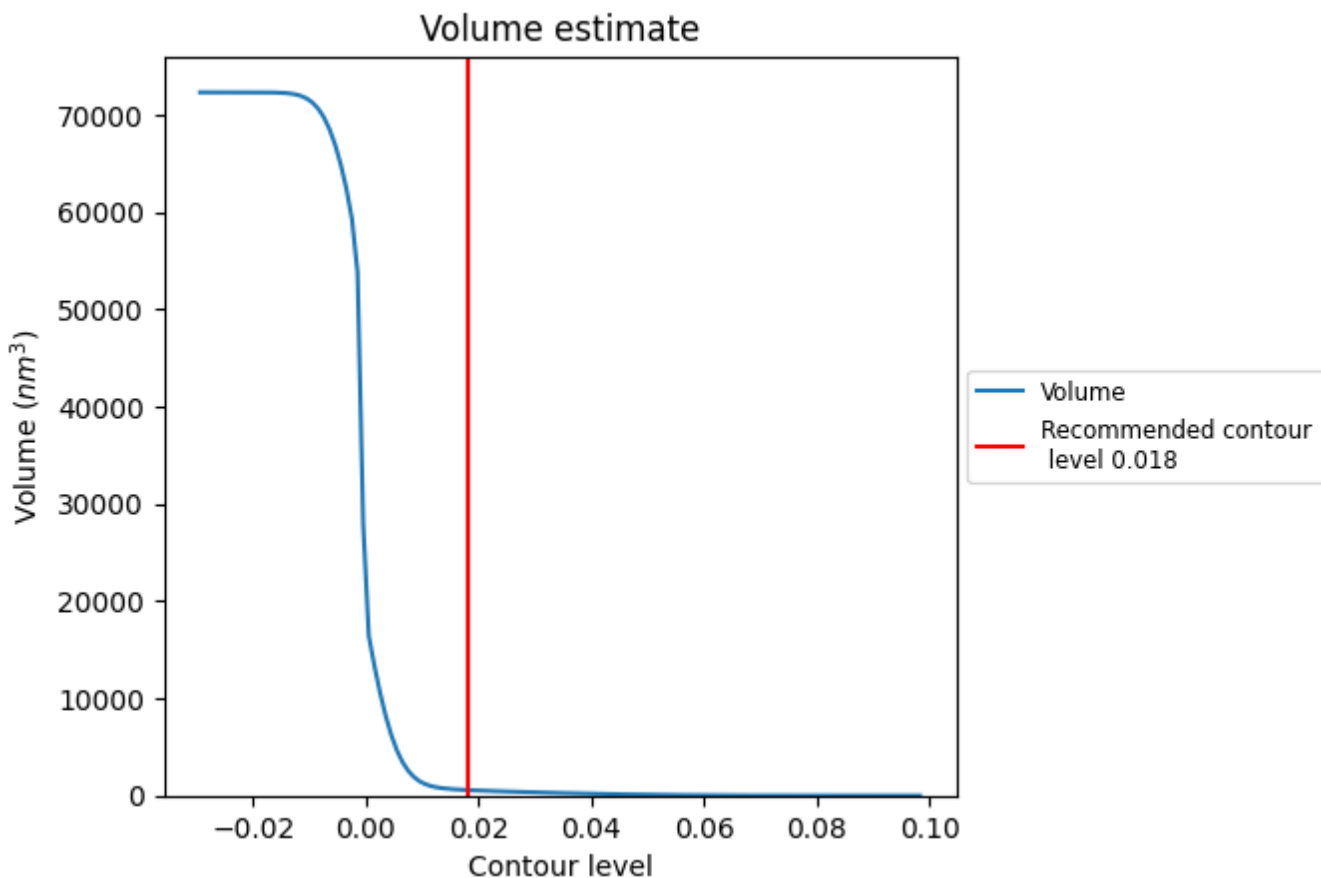
This section contains the results of statistical analysis of the map.

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.

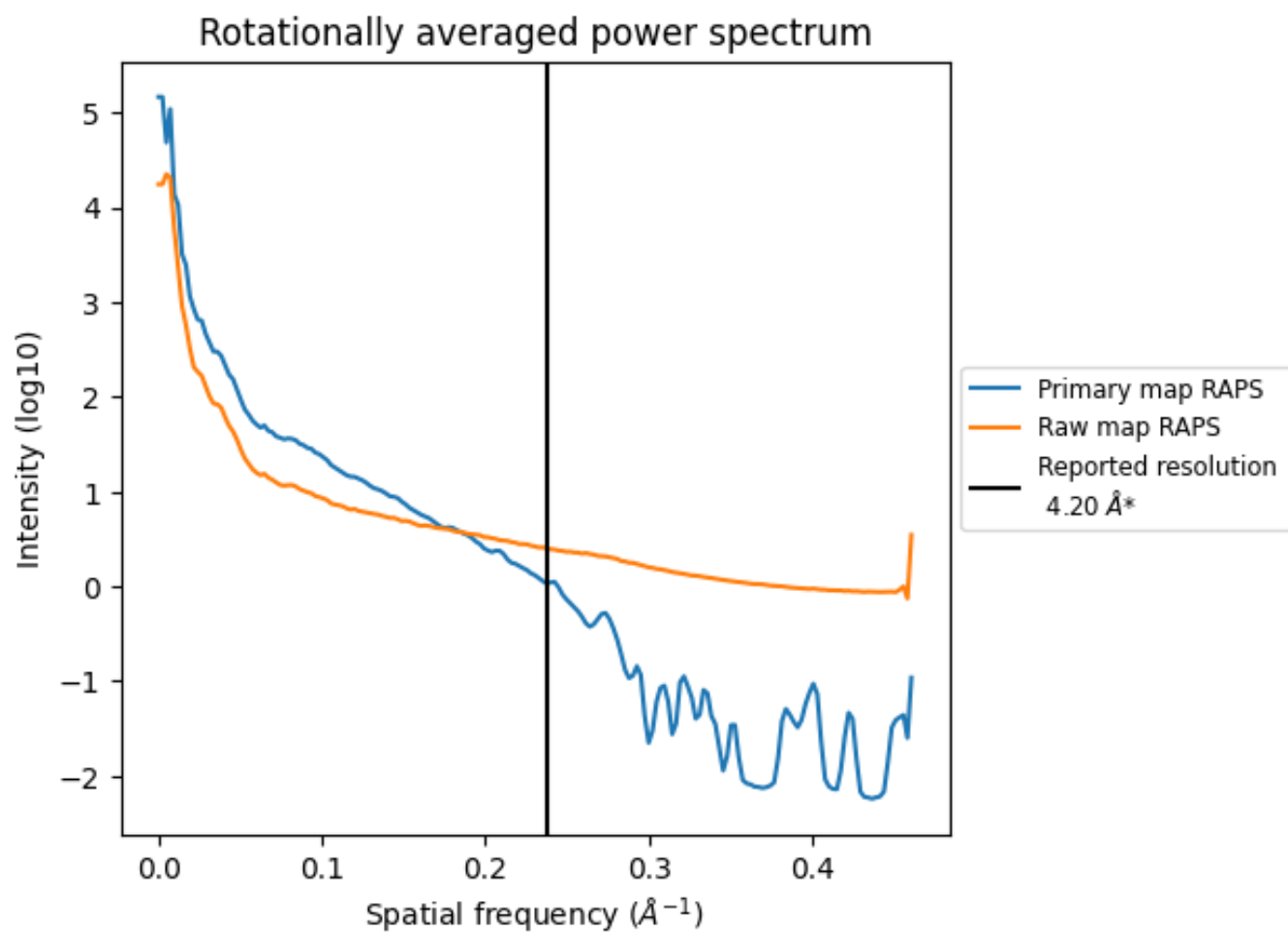
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 568 nm^3 ; this corresponds to an approximate mass of 513 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

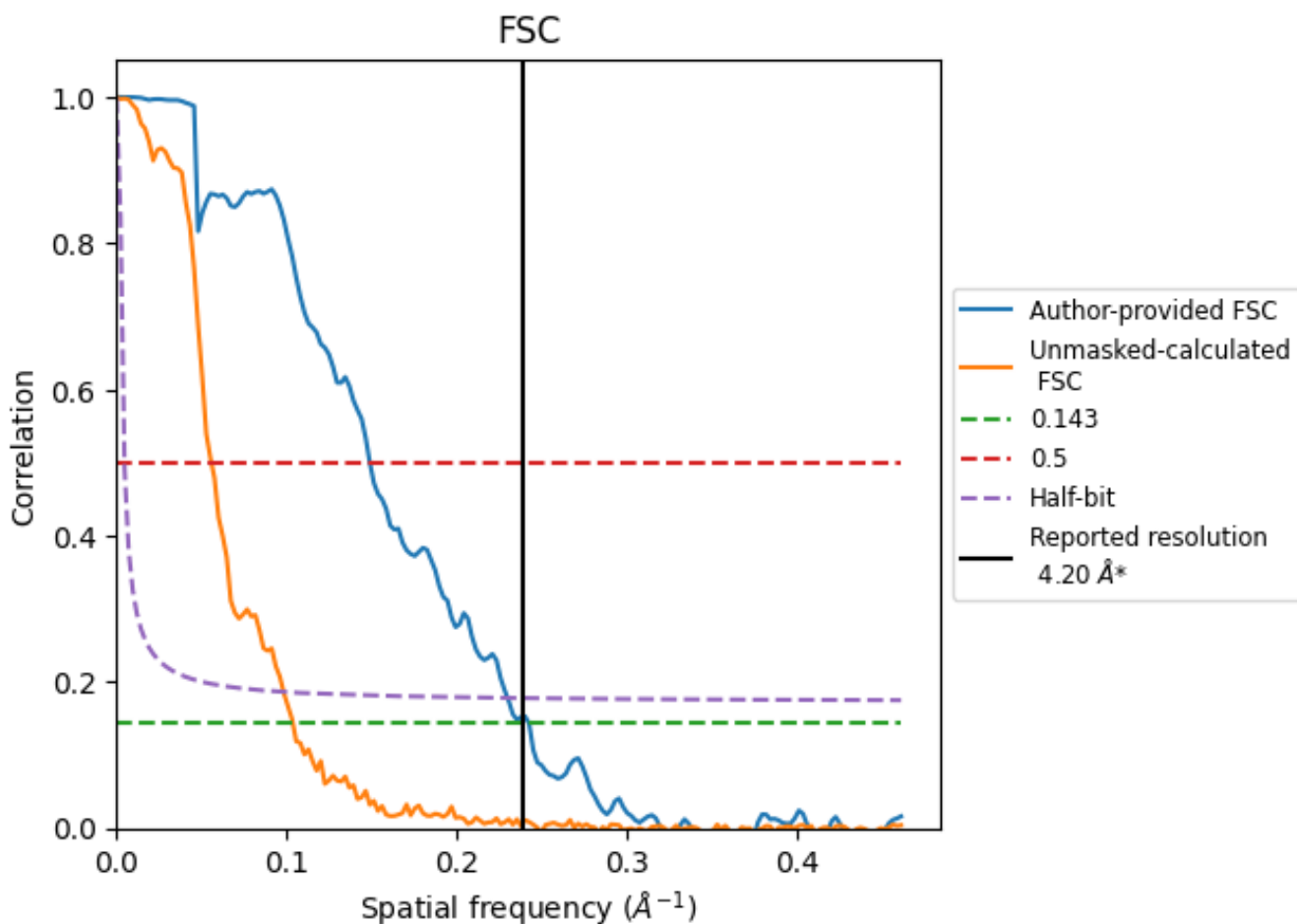


*Reported resolution corresponds to spatial frequency of 0.238 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.238 Å⁻¹

8.2 Resolution estimates [i](#)

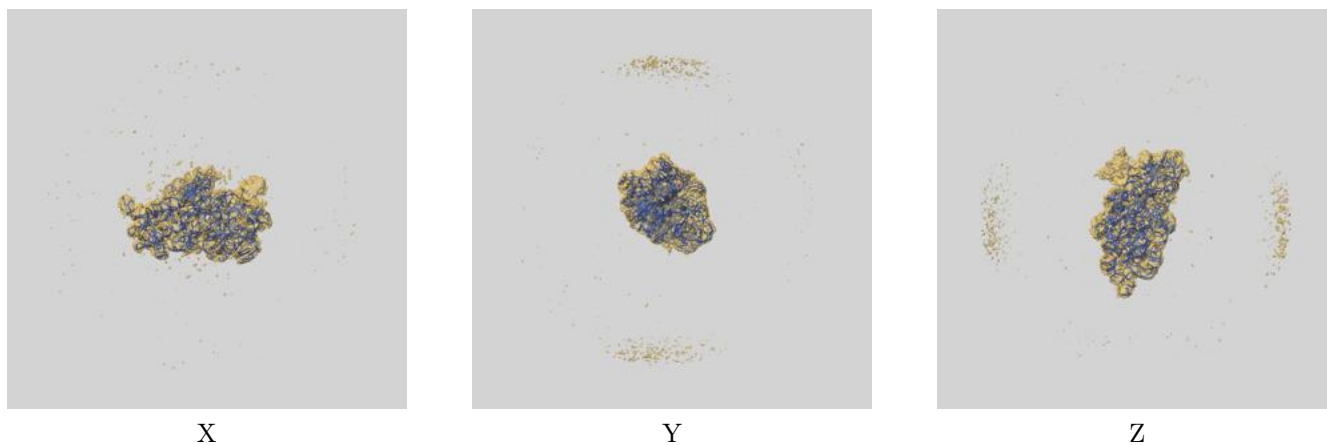
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.20	-	-
Author-provided FSC curve	4.13	6.72	4.35
Unmasked-calculated*	9.65	17.92	10.18

*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 9.65 differs from the reported value 4.2 by more than 10 %

9 Map-model fit [i](#)

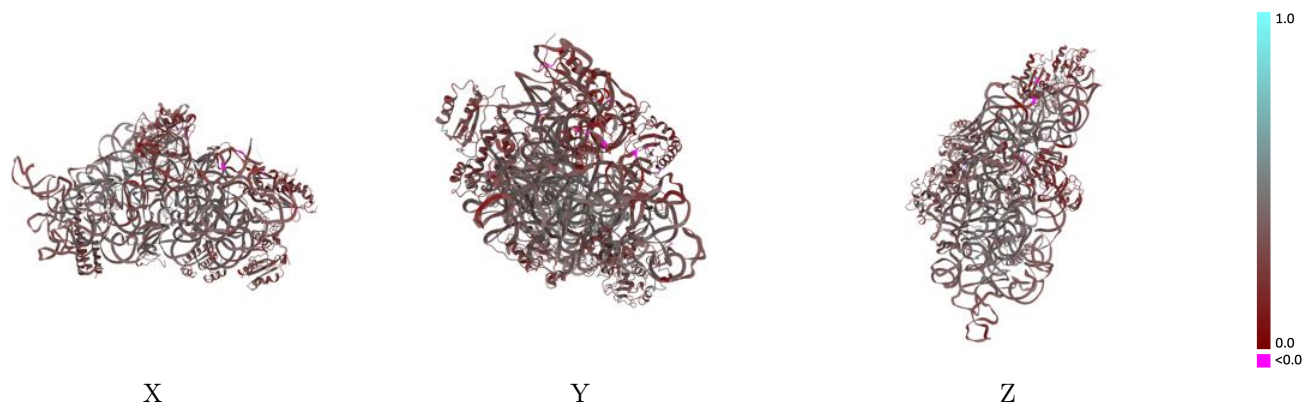
This section contains information regarding the fit between EMDB map EMD-11769 and PDB model 7AFL. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



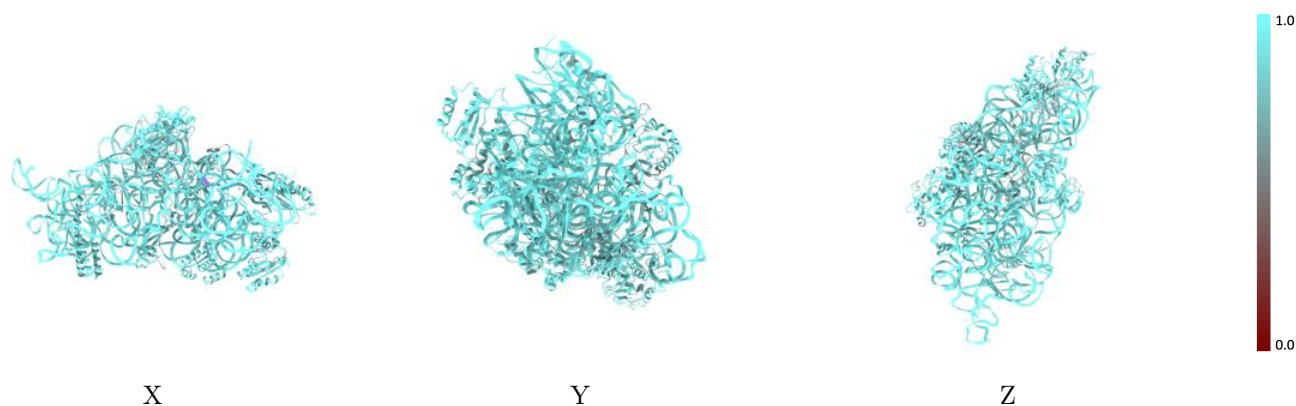
The images above show the 3D surface view of the map at the recommended contour level 0.018 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



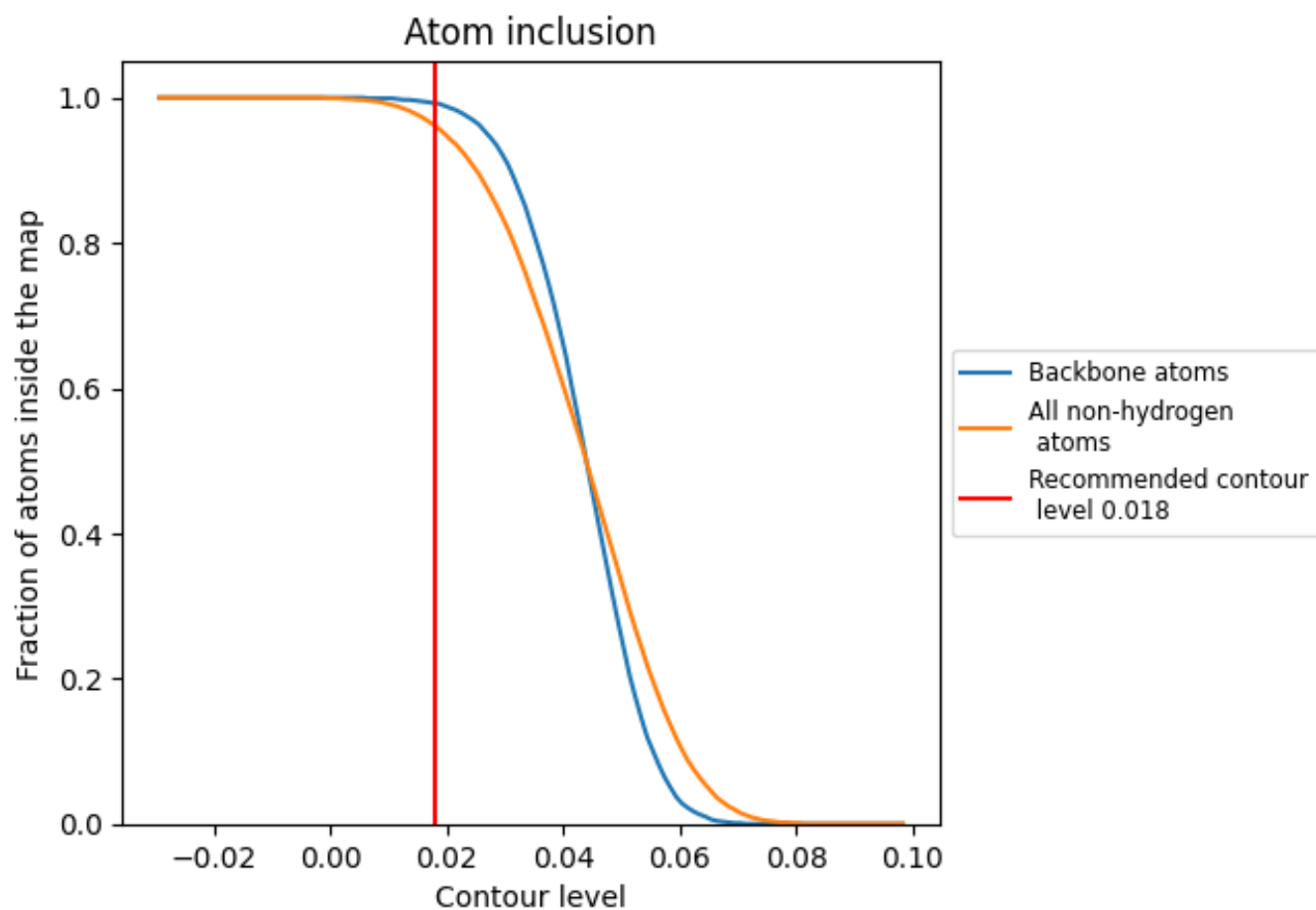
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.018).































9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 96% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.018) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9610	 0.3470
A	 0.9960	 0.3710
D	 0.8900	 0.3270
E	 0.8610	 0.3330
F	 0.9310	 0.3130
H	 0.8850	 0.3240
K	 0.8990	 0.2410
L	 0.7870	 0.2990
O	 0.9190	 0.3090
P	 0.9200	 0.3480
Q	 0.9050	 0.3560
R	 0.8930	 0.2870
T	 0.9330	 0.3170
V	 0.9110	 0.2370
X	 0.8970	 0.2270

