



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 12:15 AM UTC

PDB ID : 8PSS / pdb_00008pss
EMDB ID : EMD-17861
Title : Tilapia Lake Virus polymerase in cRNA pre-initiation state mode B (core-endo only)
Authors : Arragain, B.; Cusack, S.
Deposited on : 2023-07-13
Resolution : 2.83 Å(reported)

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:

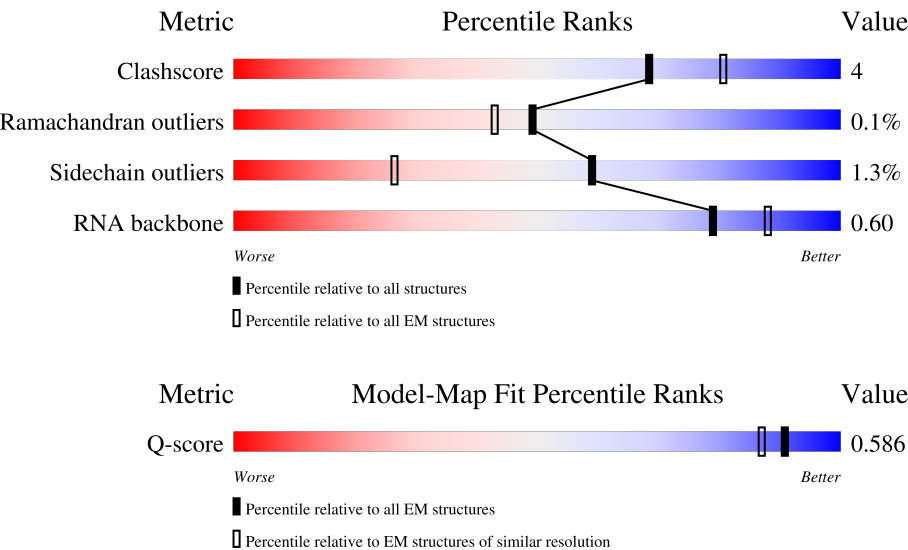
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 2.83 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.



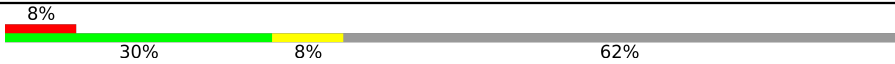
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11847 (2.33 - 3.33)

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	419	
2	B	519	
3	C	478	

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Mol	Chain	Length	Quality of chain
4	S	40	 8% 30% 8% 62%
4	V	40	 5% 22% 12% 65%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17722 atoms, of which 8690 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 Polymerase acidic protein (PA-like).

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	417	Total	C	H	N	O	S	0	0
			6652	2105	3314	597	617	19		

- Molecule 2 is a protein called Putative PB1.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	515	Total	C	H	N	O	S	0	0
			7973	2490	4000	694	759	30		

- Molecule 3 is a protein called RNA-dependent RNA polymerase.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	139	Total	C	H	N	O	S	0	0
			2221	693	1119	200	207	2		

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	391	LYS	ARG	conflict	UNP A0A7G3S745
C	458	GLY	-	expression tag	UNP A0A7G3S745
C	459	SER	-	expression tag	UNP A0A7G3S745
C	460	GLY	-	expression tag	UNP A0A7G3S745
C	461	SER	-	expression tag	UNP A0A7G3S745
C	462	GLU	-	expression tag	UNP A0A7G3S745
C	463	ASN	-	expression tag	UNP A0A7G3S745
C	464	LEU	-	expression tag	UNP A0A7G3S745
C	465	TYR	-	expression tag	UNP A0A7G3S745
C	466	PHE	-	expression tag	UNP A0A7G3S745
C	467	GLN	-	expression tag	UNP A0A7G3S745
C	468	GLY	-	expression tag	UNP A0A7G3S745
C	469	HIS	-	expression tag	UNP A0A7G3S745
C	470	HIS	-	expression tag	UNP A0A7G3S745
C	471	HIS	-	expression tag	UNP A0A7G3S745

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Chain	Residue	Modelled	Actual	Comment	Reference
C	472	HIS	-	expression tag	UNP A0A7G3S745
C	473	HIS	-	expression tag	UNP A0A7G3S745
C	474	HIS	-	expression tag	UNP A0A7G3S745
C	475	HIS	-	expression tag	UNP A0A7G3S745
C	476	HIS	-	expression tag	UNP A0A7G3S745
C	477	HIS	-	expression tag	UNP A0A7G3S745
C	478	HIS	-	expression tag	UNP A0A7G3S745

- Molecule 4 is a RNA chain called 5' cRNA end - cRNA loop (40-mer).

Mol	Chain	Residues	Atoms						AltConf	Trace
4	V	14	Total	C	H	N	O	P	0	0
			408	132	116	49	97	14		
4	S	15	Total	C	H	N	O	P	0	0
			466	145	141	61	104	15		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	A	1	Total	Zn	0
			1	1	

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
6	B	1	Total	Mg	0
			1	1	

- Molecule 4: 5' cRNA end - cRNA loop (40-mer)



- Molecule 4: 5' cRNA end - cRNA loop (40-mer)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	109188	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.089	Depositor
Minimum map value	-0.039	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	251.99998, 251.99998, 251.99998	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	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, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.10	0/3413	0.24	0/4611
2	B	0.10	0/4045	0.24	0/5474
3	C	0.10	0/1121	0.26	0/1507
4	S	0.06	0/364	0.14	0/566
4	V	0.08	0/325	0.19	0/502
All	All	0.10	0/9268	0.24	0/12660

There are no bond length outliers.

There are no bond angle outliers.

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	A	3338	3314	3313	24	0
2	B	3973	4000	4000	31	0
3	C	1102	1119	1119	10	0
4	S	325	141	162	4	0
4	V	292	116	150	2	0
5	A	1	0	0	0	0
6	B	1	0	0	0	0
All	All	9032	8690	8744	66	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 4.

All (66) 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)
2:B:199:ALA:HB1	2:B:292:ILE:HD11	1.55	0.86
4:V:13:C:O2	4:S:3:G:N2	2.14	0.80
2:B:335:THR:OG1	2:B:337:ASP:OD1	1.99	0.79
1:A:235:THR:HG21	1:A:305:SER:HB2	1.66	0.78
1:A:45:TYR:OH	1:A:50:ASN:OD1	2.04	0.74
1:A:81:CYS:SG	1:A:82:SER:N	2.62	0.72
2:B:103:THR:HG21	2:B:111:LEU:HD13	1.73	0.69
3:C:1:MET:HG3	3:C:6:LYS:HE3	1.75	0.67
2:B:381:GLU:OE2	2:B:419:SER:OG	2.15	0.65
3:C:62:ARG:NH1	3:C:63:PRO:O	2.30	0.65
1:A:58:GLU:OE2	1:A:74:ASN:ND2	2.35	0.60
1:A:9:THR:HG21	1:A:43:ASP:HB2	1.83	0.59
1:A:287:GLU:OE1	1:A:292:ARG:NH1	2.36	0.59
1:A:30:SER:OG	1:A:33:GLU:OE1	2.11	0.58
2:B:145:ARG:NH2	4:S:13:U:O2	2.35	0.57
3:C:89:GLU:OE1	3:C:89:GLU:N	2.34	0.57
2:B:219:GLU:N	2:B:219:GLU:OE2	2.38	0.56
2:B:466:SER:O	2:B:487:ARG:NH2	2.40	0.55
1:A:218:PRO:HD2	1:A:348:VAL:HG11	1.89	0.54
2:B:73:ARG:NH1	4:S:14:G:OP2	2.41	0.54
3:C:1:MET:CG	3:C:6:LYS:HE3	2.38	0.54
1:A:19:SER:OG	1:A:20:GLU:OE1	2.26	0.53
2:B:199:ALA:HB1	2:B:292:ILE:CD1	2.34	0.52
1:A:26:LEU:O	1:A:26:LEU:HD12	2.09	0.51
1:A:235:THR:HG21	1:A:305:SER:CB	2.39	0.51
3:C:78:SER:OG	3:C:114:ASN:OD1	2.27	0.51
2:B:467:ASP:O	2:B:467:ASP:OD1	2.29	0.50
2:B:35:SER:O	2:B:39:THR:HG23	2.12	0.48
4:V:13:C:N3	4:S:3:G:N1	2.49	0.48
2:B:65:VAL:HG23	2:B:65:VAL:O	2.14	0.47
2:B:418:ILE:HD11	2:B:428:CYS:HB3	1.97	0.47
1:A:301:ASP:C	1:A:301:ASP:OD1	2.58	0.46
2:B:277:LEU:HD22	2:B:313:HIS:CD2	2.50	0.46
1:A:244:SER:OG	1:A:245:ASP:N	2.49	0.46
1:A:23:ARG:NH1	1:A:33:GLU:OE2	2.49	0.46
2:B:23:PRO:HA	2:B:144:LEU:HD13	1.98	0.46
2:B:330:ASN:O	2:B:331:SER:OG	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:ASP:O	1:A:50:ASN:N	2.50	0.45
2:B:103:THR:HG23	2:B:125:ILE:HG22	1.97	0.45
2:B:111:LEU:HD12	2:B:111:LEU:O	2.17	0.45
3:C:1:MET:SD	3:C:2:SER:N	2.90	0.44
3:C:81:SER:OG	3:C:113:VAL:HG13	2.17	0.44
1:A:170:GLU:HB3	1:A:223:LEU:HD11	1.99	0.44
1:A:55:PHE:CD1	1:A:55:PHE:N	2.86	0.44
1:A:176:GLU:H	1:A:176:GLU:CD	2.26	0.44
2:B:134:LEU:HD21	2:B:241:MET:HG2	2.00	0.43
2:B:418:ILE:HD11	2:B:428:CYS:CB	2.48	0.43
2:B:24:ASP:OD1	2:B:24:ASP:N	2.49	0.43
2:B:467:ASP:OD1	2:B:467:ASP:C	2.61	0.43
2:B:171:VAL:HG21	2:B:242:LEU:HD12	2.00	0.42
1:A:133:ARG:HH11	1:A:133:ARG:HG3	1.84	0.42
1:A:163:LYS:O	1:A:296:HIS:CG	2.72	0.42
2:B:103:THR:O	2:B:103:THR:HG22	2.19	0.42
2:B:458:ARG:NH2	3:C:24:VAL:O	2.53	0.42
1:A:396:ILE:HB	1:A:397:PRO:HD3	2.00	0.42
3:C:55:ASP:N	3:C:55:ASP:OD1	2.53	0.42
2:B:277:LEU:HD23	2:B:286:SER:CB	2.50	0.42
2:B:165:ARG:O	2:B:169:SER:OG	2.36	0.41
1:A:1:MET:SD	1:A:2:ASP:N	2.94	0.41
1:A:16:PHE:CD2	1:A:89:ILE:HD11	2.54	0.41
2:B:42:SER:HG	2:B:43:TYR:HD2	1.64	0.41
2:B:504:ASN:OD1	2:B:505:PHE:N	2.54	0.41
2:B:482:ASP:O	2:B:486:ILE:HG13	2.20	0.41
2:B:481:SER:O	2:B:485:LYS:HG3	2.21	0.41
1:A:172:GLY:O	1:A:173:LEU:HD12	2.20	0.41
3:C:124:ASN:OD1	3:C:124:ASN:N	2.53	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	A	415/419 (99%)	404 (97%)	11 (3%)	0	100	100
2	B	513/519 (99%)	502 (98%)	10 (2%)	1 (0%)	43	62
3	C	137/478 (29%)	132 (96%)	5 (4%)	0	100	100
All	All	1065/1416 (75%)	1038 (98%)	26 (2%)	1 (0%)	49	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	254	LYS

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	A	364/365 (100%)	363 (100%)	1 (0%)	86	94
2	B	445/447 (100%)	438 (98%)	7 (2%)	55	76
3	C	121/405 (30%)	117 (97%)	4 (3%)	33	58
All	All	930/1217 (76%)	918 (99%)	12 (1%)	59	79

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	SER
2	B	59	GLU
2	B	71	SER
2	B	75	CYS
2	B	285	LEU
2	B	401	THR
2	B	444	ASP
2	B	468	LEU
3	C	3	GLN
3	C	50	GLN
3	C	67	SER
3	C	125	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	74	ASN
1	A	87	GLN
1	A	246	ASN
1	A	369	ASN
2	B	296	ASN
2	B	385	GLN
3	C	3	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	S	14/40 (35%)	0	0
4	V	13/40 (32%)	4 (30%)	0
All	All	27/80 (33%)	4 (14%)	0

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	V	5	A
4	V	7	U
4	V	10	A
4	V	11	C

There are no RNA pucker outliers to report.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 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.

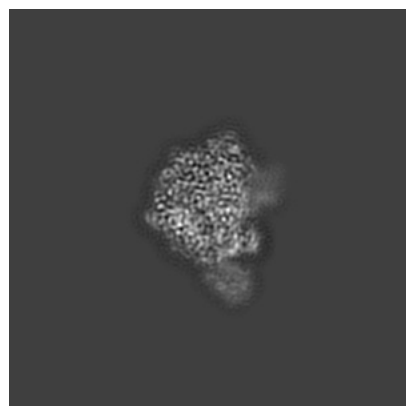
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17861. These allow visual inspection of the internal detail of the map and identification of artifacts.

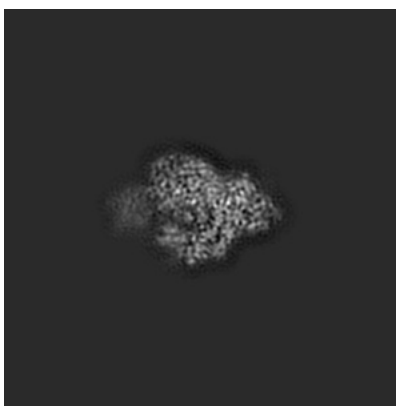
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

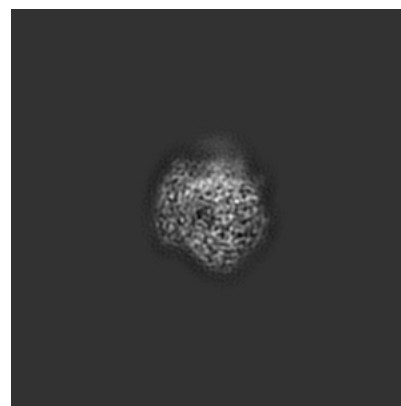
6.1.1 Primary map



X

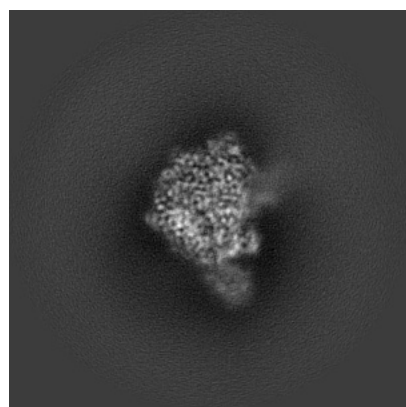


Y

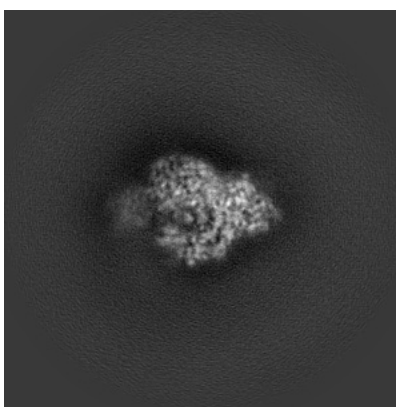


Z

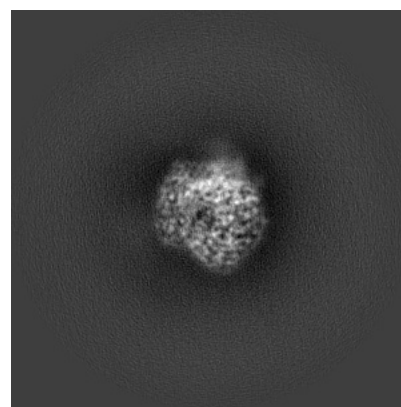
6.1.2 Raw map



X



Y

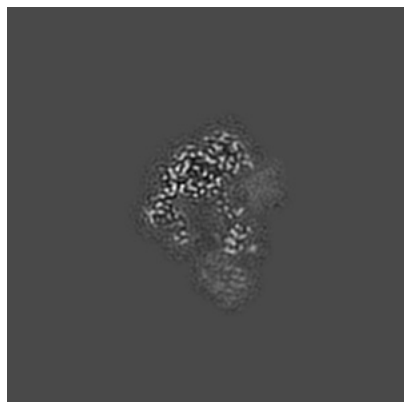


Z

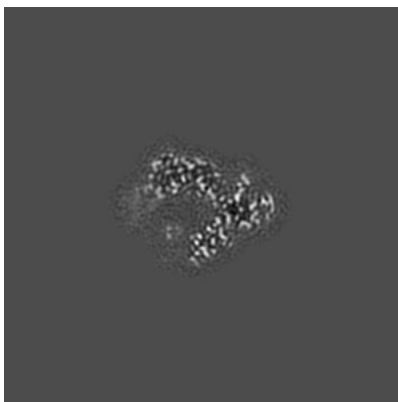
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

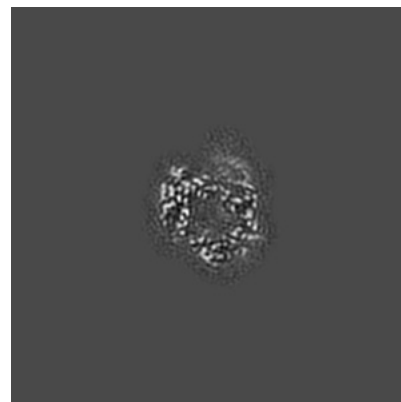
6.2.1 Primary map



X Index: 150

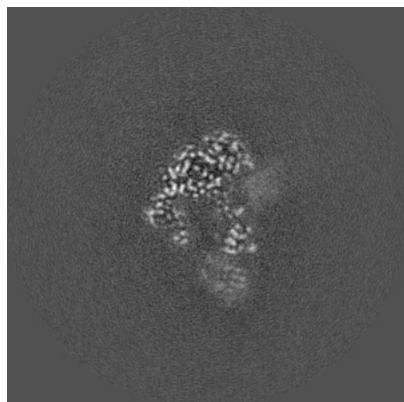


Y Index: 150

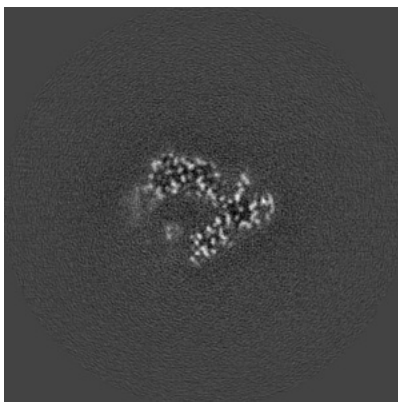


Z Index: 150

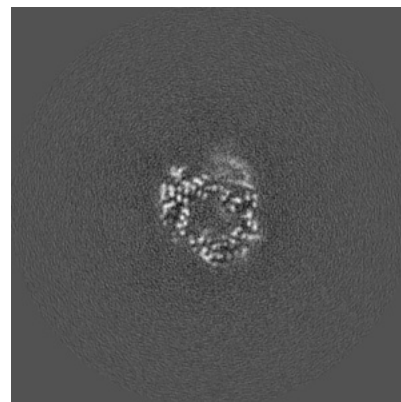
6.2.2 Raw map



X Index: 150



Y Index: 150

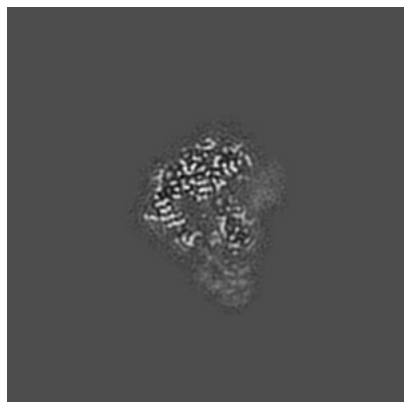


Z Index: 150

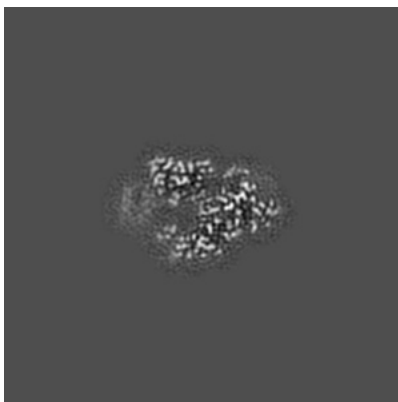
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 155

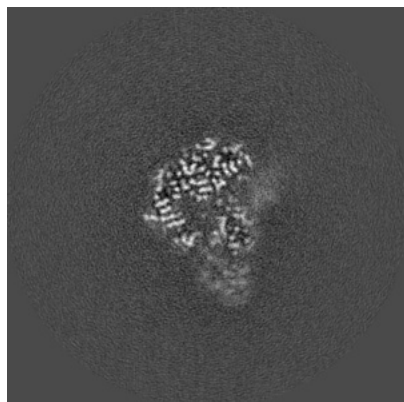


Y Index: 158

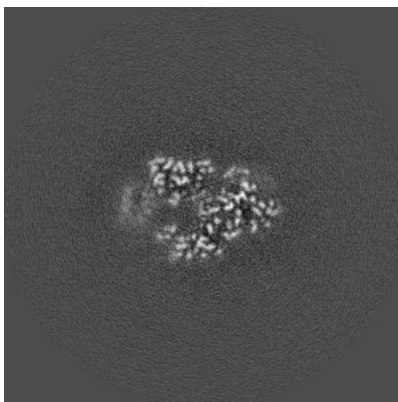


Z Index: 148

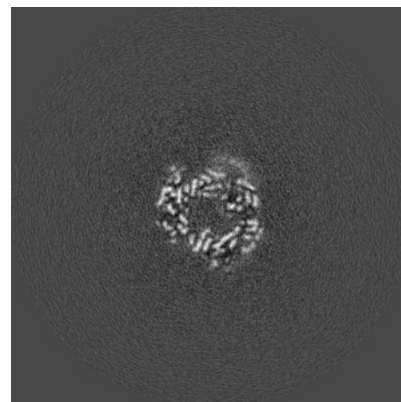
6.3.2 Raw map



X Index: 155



Y Index: 158

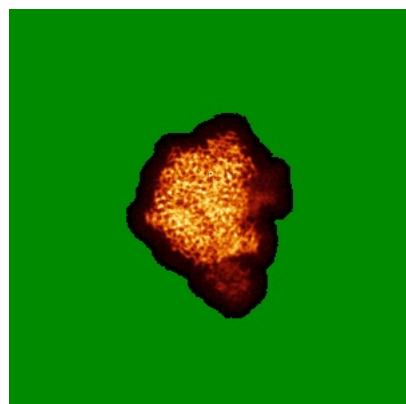


Z Index: 148

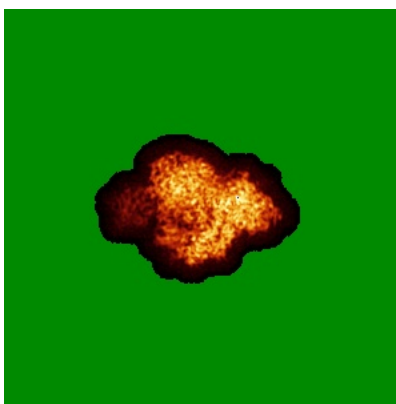
The images above show the largest variance slices of the map in three orthogonal directions.

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

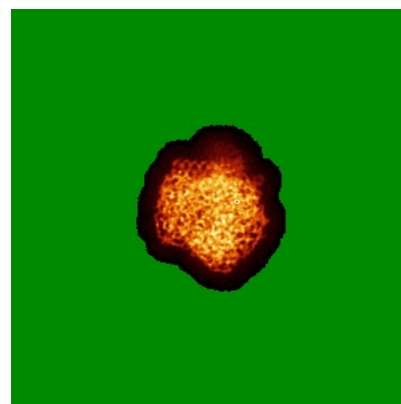
6.4.1 Primary map



X

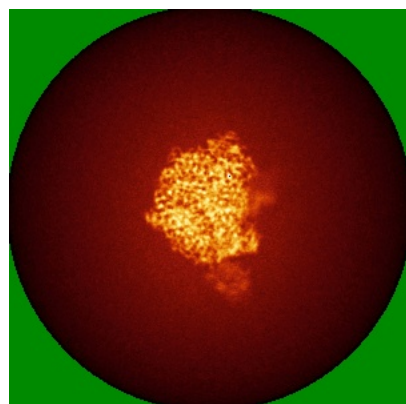


Y

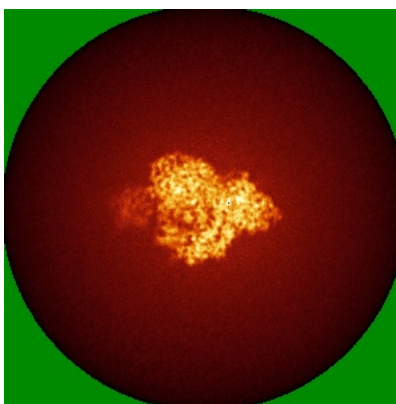


Z

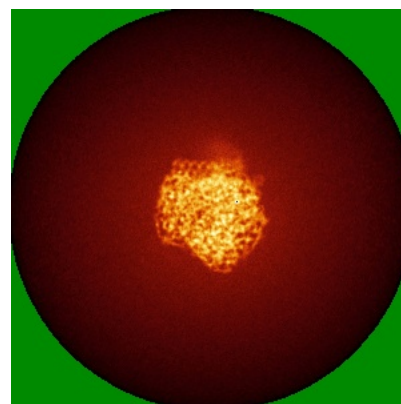
6.4.2 Raw map



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](#)

This section was not generated.

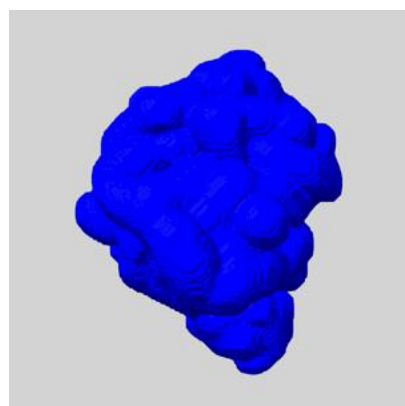
6.6 Mask visualisation [i](#)

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

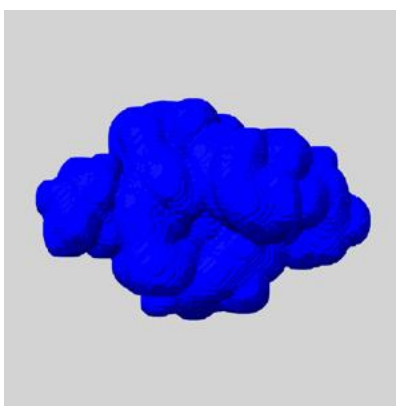
A mask typically either:

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

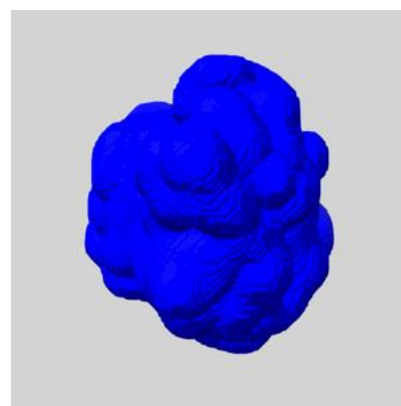
6.6.1 emd_17861_msk_1.map [i](#)



X



Y

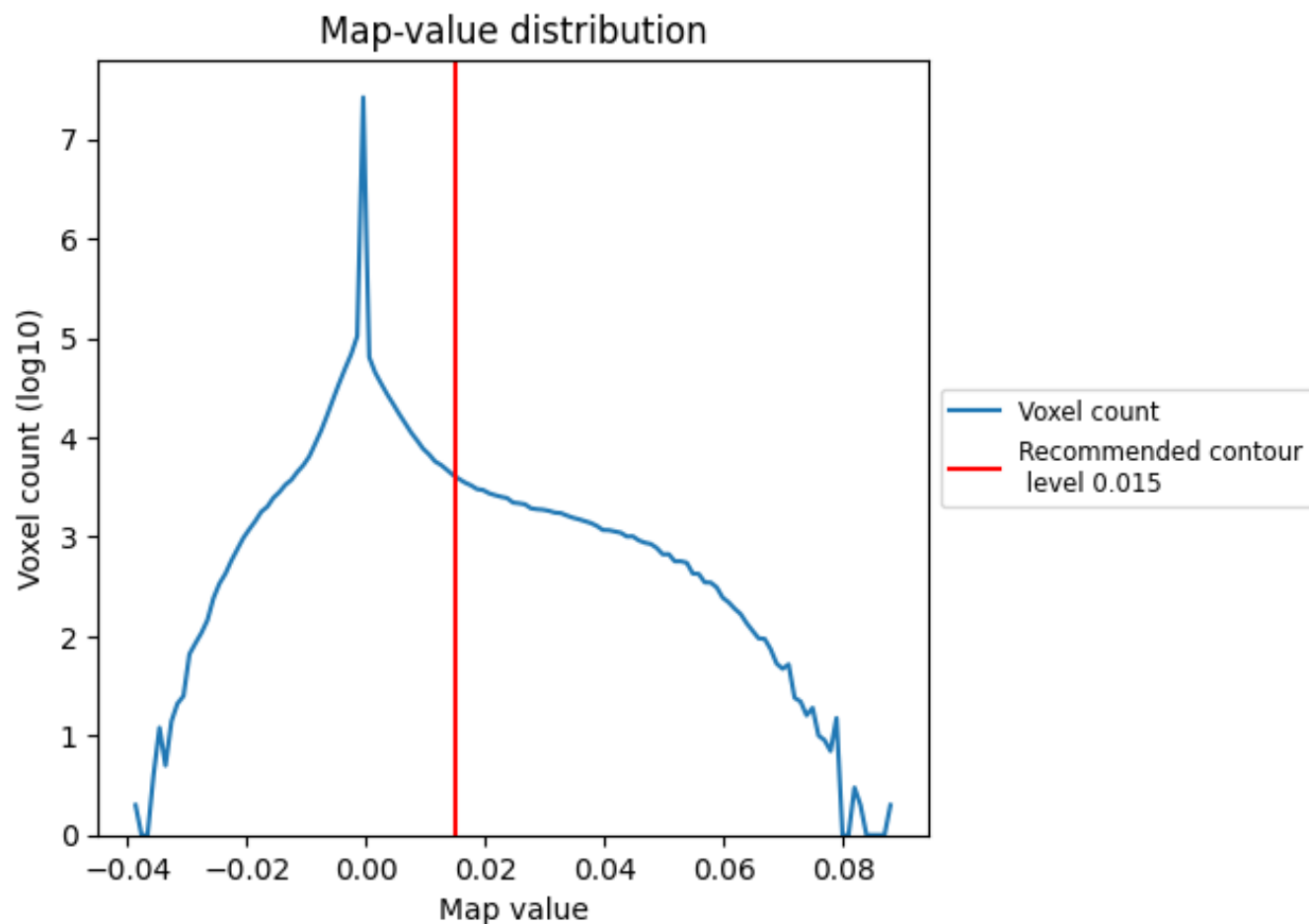


Z

7 Map analysis [i](#)

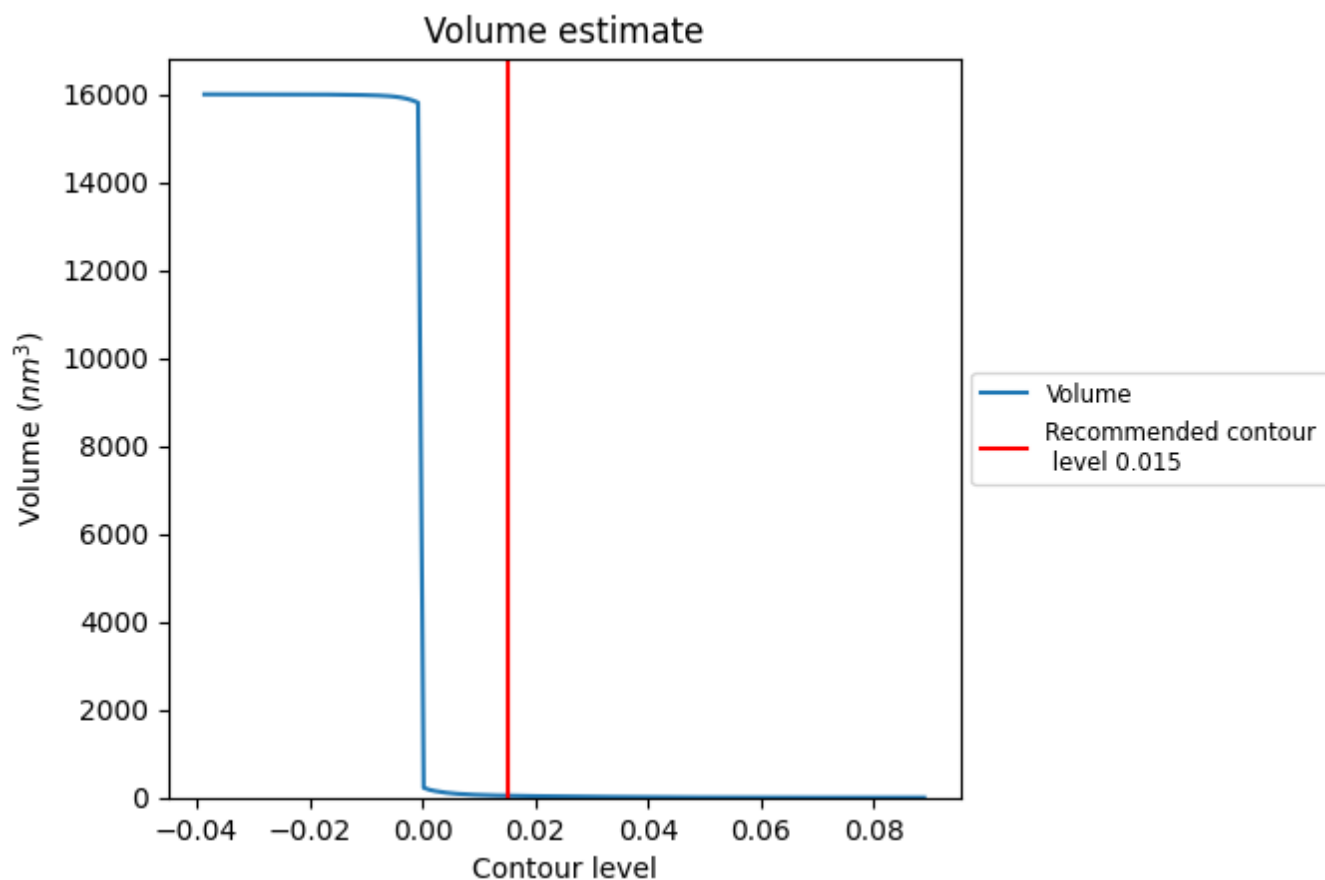
This section contains the results of statistical analysis of the map.

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.

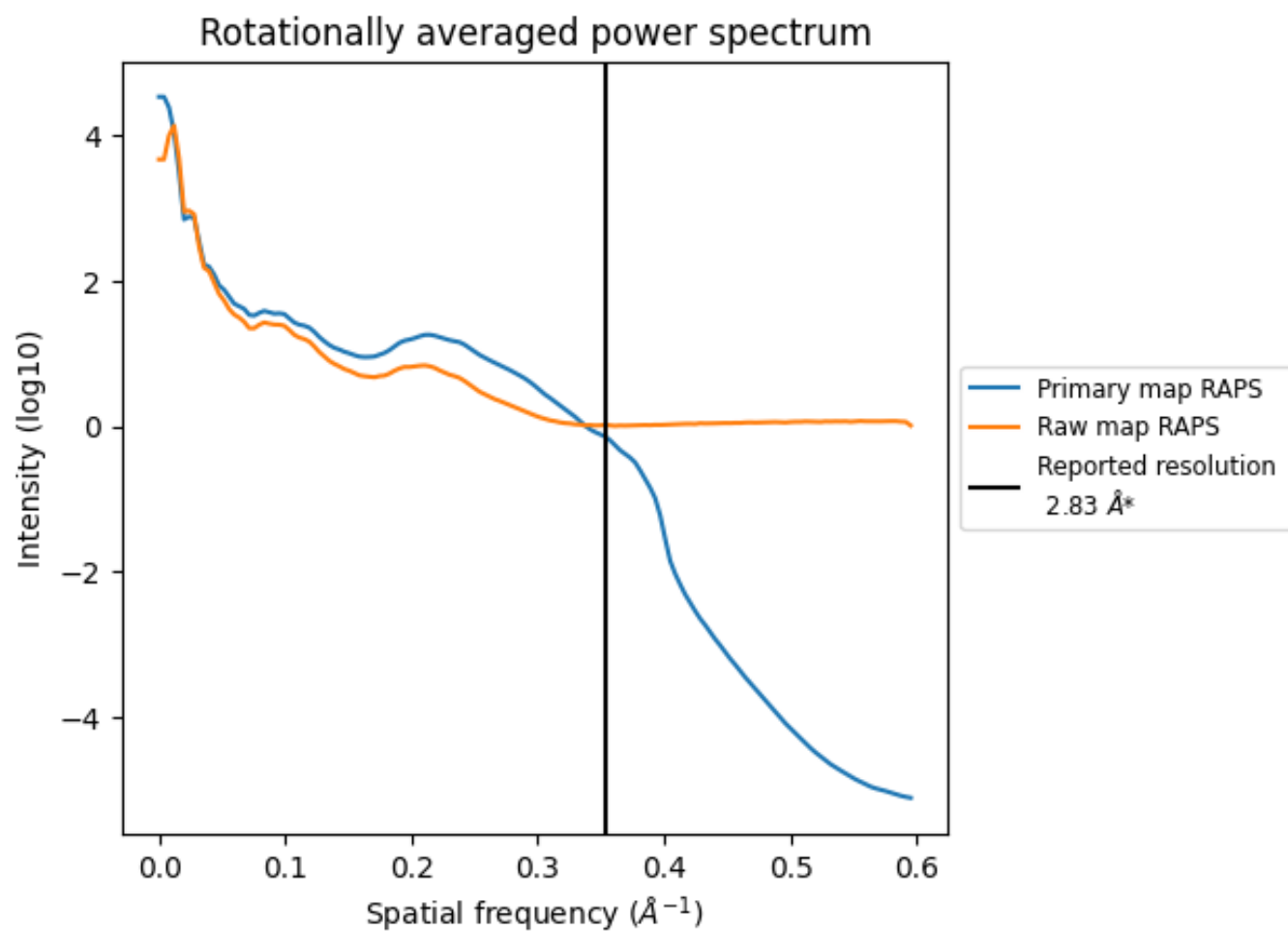
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 43 nm^3 ; this corresponds to an approximate mass of 39 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 ⓘ

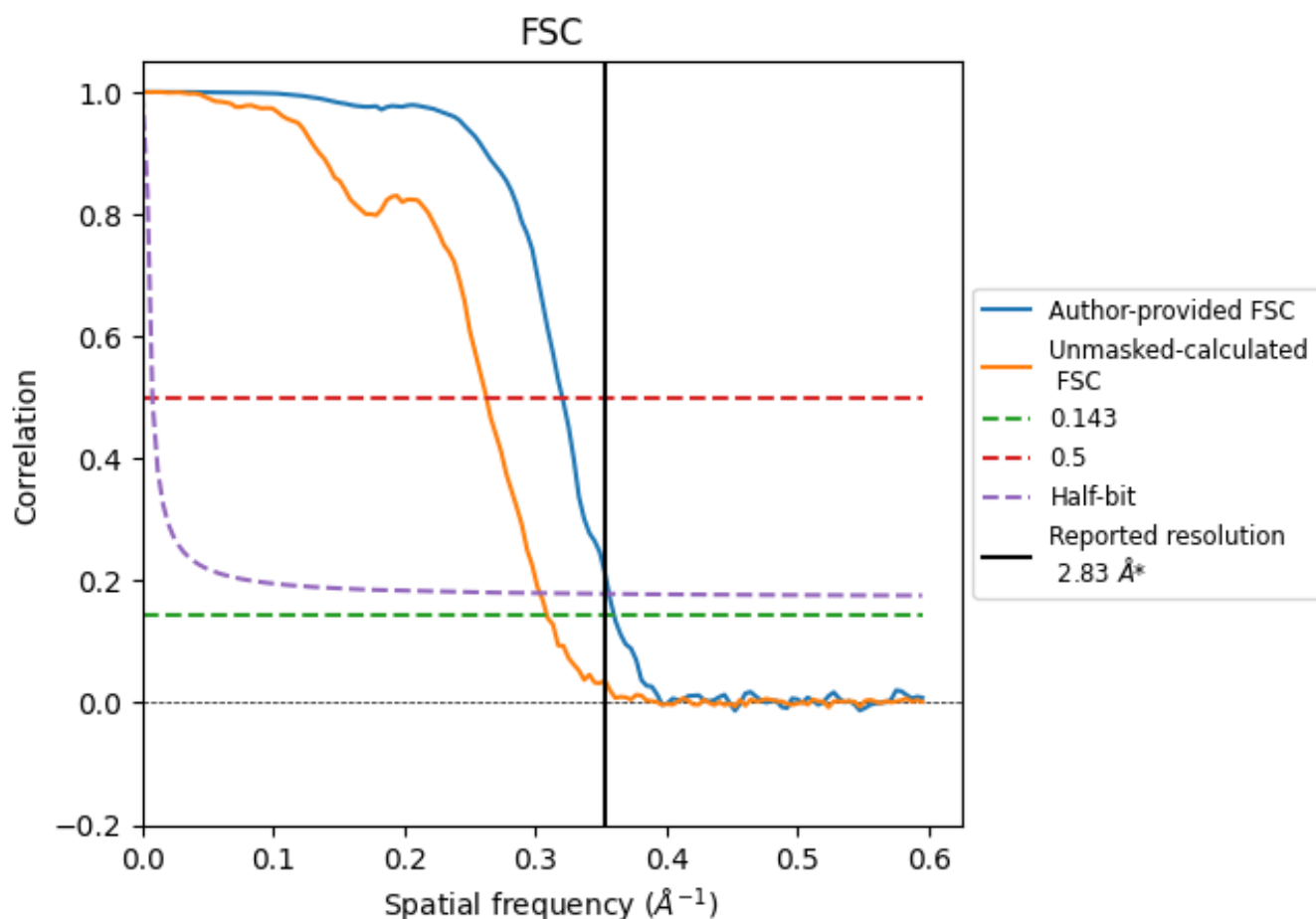


*Reported resolution corresponds to spatial frequency of 0.353 Å⁻¹

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 0.353 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.83	-	-
Author-provided FSC curve	2.78	3.12	2.81
Unmasked-calculated*	3.24	3.81	3.29

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.24 differs from the reported value 2.83 by more than 10 %

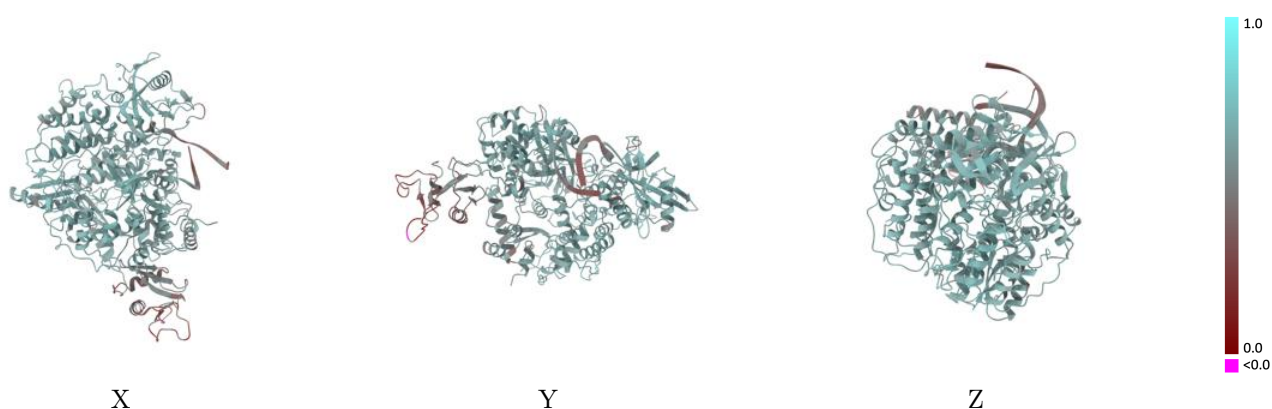
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-17861 and PDB model 8PSS. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

This section was not generated.

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

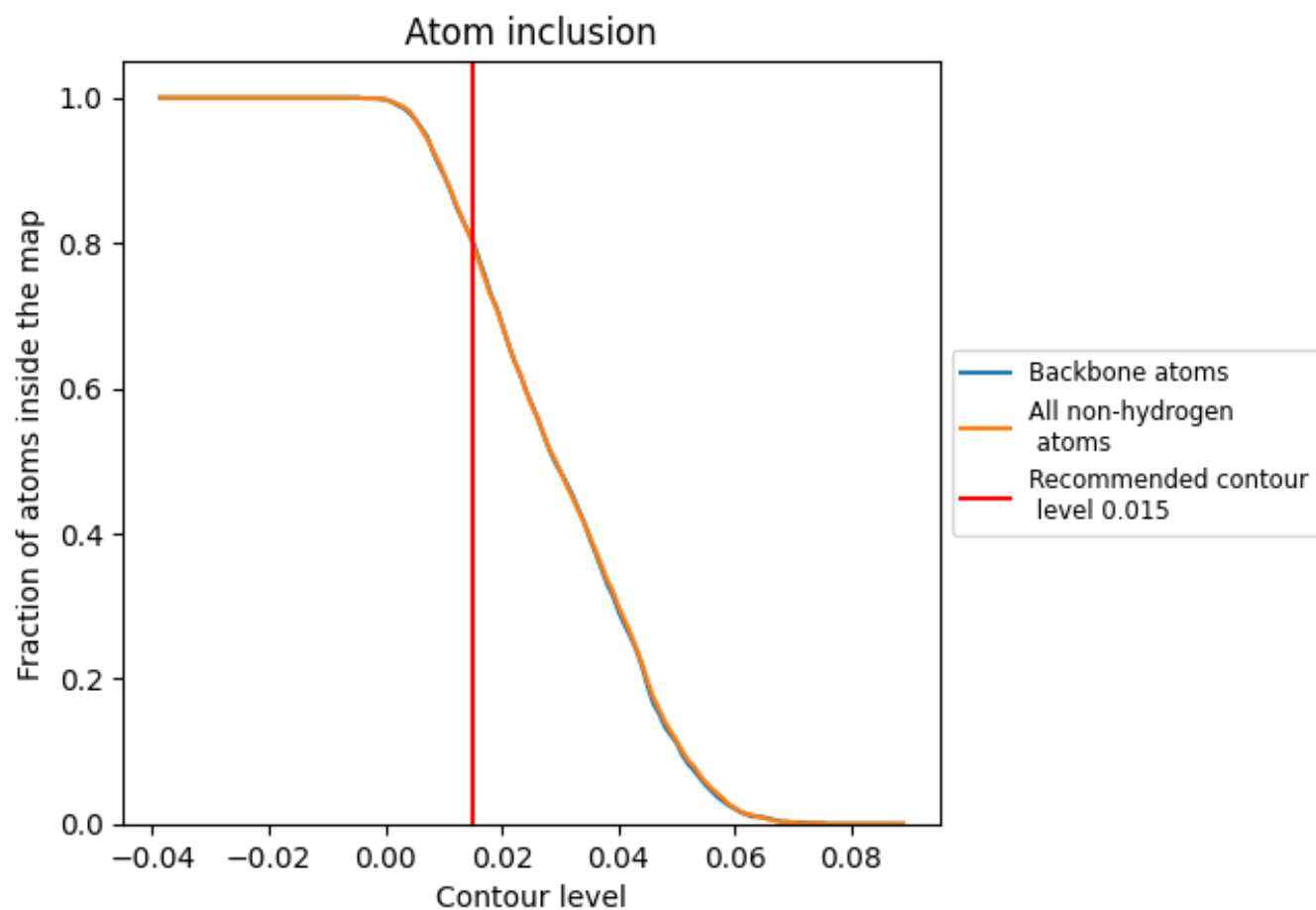


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 80% 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 (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8010	0.5860
A	0.7170	0.5560
B	0.8790	0.6160
C	0.8300	0.6010
S	0.6770	0.5320
V	0.7160	0.5060

