



## wwPDB EM Validation Summary Report ⓘ

Mar 6, 2026 – 06:48 AM UTC

PDB ID : 8PSO / pdb\_00008pso  
EMDB ID : EMD-17858  
Title : Tilapia Lake Virus polymerase in vRNA initiation state (core only)  
Authors : Arragain, B.; Cusack, S.  
Deposited on : 2023-07-13  
Resolution : 2.40 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:

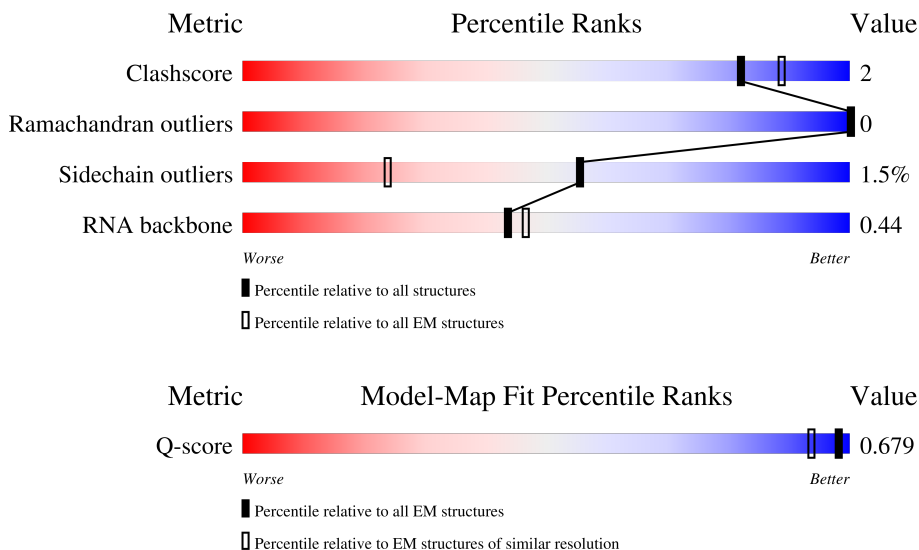
EMDB validation analysis : 0.0.1.dev132  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)  
MapQ : 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 2.40 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.





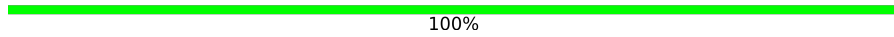
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	5628 ( 1.90 - 2.90 )

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	A	419	
2	B	519	
3	C	478	

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Mol	Chain	Length	Quality of chain
4	S	40	
4	V	40	
5	F	1	

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 16213 atoms, of which 7949 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	
			Total	C	H	N	O			S
1	A	316	5036	1584	2524	459	457	12	0	0

- Molecule 2 is a protein called Putative PB1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	B	515	7974	2490	4001	694	759	30	1	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
3	C	140	2233	697	1123	201	210	2	0	0

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' vRNA end - vRNA loop (40-mer).

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
4	V	13	Total 406	C 121	H 138	N 42	O 92	P 13	0	0
4	S	15	Total 476	C 145	H 151	N 60	O 105	P 15	0	0

- Molecule 5 is a DNA chain called DNA (5'-D\*(CTP))-3').

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
5	F	1	Total 41	C 9	H 12	N 3	O 14	P 3	0	0

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
7	B	2	Total 2	Mg 2	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		AltConf
8	A	10	Total 10	O 10	0

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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>AltConf</b>
8	B	27	Total 27	O 27	0
8	V	4	Total 4	O 4	0
8	S	2	Total 2	O 2	0
8	F	1	Total 1	O 1	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: Polymerase acidic protein (PA-like)

Chain A:  73% 25%

MET ASP SER ARG PHE PRO PHE ALA GLN LEU THR ARG GLY VAL PHE CYS ASP ASP PHE THR TVR SER GLU GLY SER ARG ARG ARG PHE LEU SER SER TVR THR VAL GLN VAL GLN ARG ARG PRO GLY VAL PRO VAL GLN ASP GLY ASP CYS TYR ASP CYS LEU LYS ASN LYS TRP LYS ILE ALA PHE LEU LEU LEU GLN

PRO ARG LYS PHE PRO LYS ALA THR VAL ARG CYS ILE LEU ASN ASP ASP ALA THR TVR VAL CYS VAL GLN SER GLU GLN GLN TTR GLN SER GLN ILE CYS LYS VAL GLN PHE LYS ASP PRO TVR LEU LEU ILE ASP G102 V103 S111 I134 R133 E135 D144 T235 D245 G259 R260 H261 S262

R263 R264 R265 S305 Y320 D323 D337 G341 S404 F417 ALA ARG

- Molecule 2: Putative PB1

Chain B:  7% 91% 8%

M1 E6 L13 R63 I64 C75 V78 K81 E82 G83 R84 V85 K114 V118 E119 I125 T128 L134 L166 V171 S177 S183 A184 T187 A188 T189 A190 S191 A192 E193 A199 N211 D222 M241 L242 R247 M266

T281 D282 L285 S288 I292 N330 S331 D371 D406 I418 S419 F420 L424 D444 R461 L462 E463 A464 L465 S466 P470 L471 R475 S479 D482 I486 R487 A488 Q491 A492 H493 L494 K495 S496 L497 G498 L499 E500 Q501 P502 T503 F505 N506

Y507 Y510 V513 Q514 P515 THR ALA GLY CYS

- Molecule 3: RNA-dependent RNA polymerase

Chain C:  27% 71%

M1 S2 T17 S30 S67 S78 S81 F82 R83 T87 Q88 E89 E90 D93 L94 N96 N97 D100 G101 D102 R111 V112 V113 M114 A122 A139 D140 ARG ALA PHE ASP THR CYS LEU SER GLY PHE VAL ARG ALA PRO ILE LEU THR THR PRO CYS LEU PHE

ILE CYS CYS GLY PRO PRO SER PHE LYS ASP SER LEU PHE VAL ILE LYS GLY GLU PHE TRP HIS MET TYR ASP GLY PHE GLN HIS GLN PHE VAL ALA VAL VAL ASP M114 LYS PHE LEU ALA LYS SER SER PRO SER PHE THR TRP LEU ALA LYS ARG VAL ALA ARG LEU ASN LEU VAL



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	271170	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	0.223	Depositor
Minimum map value	-0.097	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.035	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: ZN, MG, CTP

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	0.09	0/2568	0.25	0/3470
2	B	0.11	0/4050	0.27	1/5481 (0.0%)
3	C	0.09	0/1129	0.25	0/1518
4	S	0.09	0/364	0.26	0/566
4	V	0.07	0/297	0.22	0/458
All	All	0.10	0/8408	0.26	1/11493 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	470	PRO	CA-N-CD	-6.39	103.06	112.00

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	2512	2524	2523	10	0
2	B	3973	4001	3995	24	0
3	C	1110	1123	1123	8	0
4	S	325	151	161	4	0
4	V	268	138	139	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	29	12	12	0	0
6	A	1	0	0	0	0
7	B	2	0	0	0	0
8	A	10	0	0	1	0
8	B	27	0	0	0	0
8	F	1	0	0	0	0
8	S	2	0	0	0	0
8	V	4	0	0	0	0
All	All	8264	7949	7953	40	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

The worst 5 of 40 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)
4:V:10:U:O4	4:S:10:A:N6	2.17	0.77
3:C:78:SER:OG	3:C:114:ASN:OD1	2.07	0.71
2:B:510:TYR:OH	3:C:90:GLU:OE2	2.09	0.70
1:A:404:SER:OG	8:A:601:HOH:O	2.13	0.66
2:B:444:ASP:OD1	2:B:444:ASP:N	2.31	0.63

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	
1	A	314/419 (75%)	309 (98%)	5 (2%)	0	100	100
2	B	514/519 (99%)	503 (98%)	11 (2%)	0	100	100
3	C	138/478 (29%)	134 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	966/1416 (68%)	946 (98%)	20 (2%)	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	
1	A	272/365 (74%)	271 (100%)	1 (0%)	84	92
2	B	445/447 (100%)	437 (98%)	8 (2%)	51	73
3	C	122/405 (30%)	118 (97%)	4 (3%)	33	55
All	All	839/1217 (69%)	826 (98%)	13 (2%)	55	77

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	444	ASP
2	B	486	ILE
3	C	87	THR
3	C	30	SER
3	C	67	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	421	GLN
2	B	514	GLN
3	C	128	HIS
2	B	313	HIS
2	B	211	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	S	15/40 (37%)	3 (20%)	1 (6%)
4	V	12/40 (30%)	4 (33%)	0
All	All	27/80 (33%)	7 (25%)	1 (3%)

5 of 7 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	V	5	A
4	V	7	C
4	V	10	U
4	V	11	C
4	S	2	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	S	1	G

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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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.

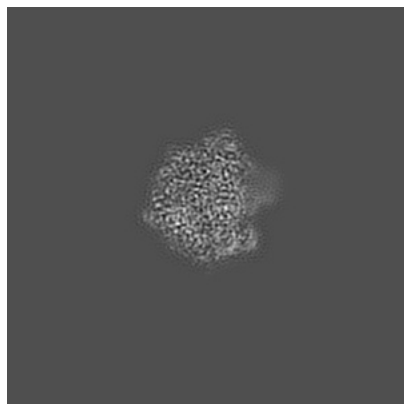
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17858. These allow visual inspection of the internal detail of the map and identification of artifacts.

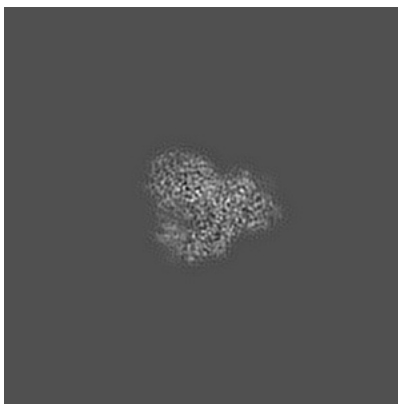
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

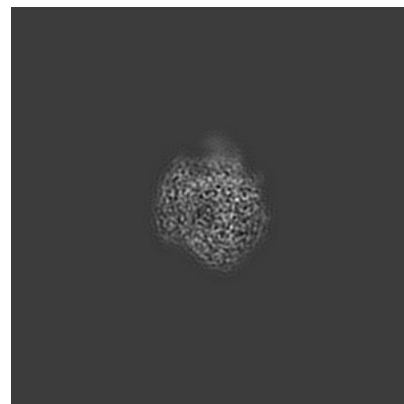
#### 6.1.1 Primary map



X

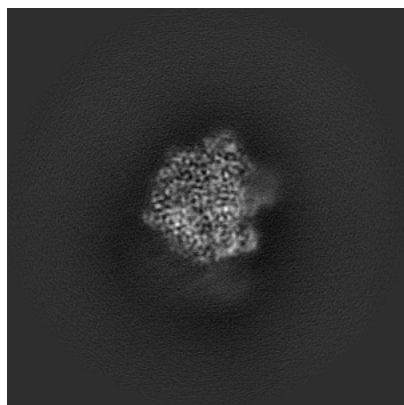


Y

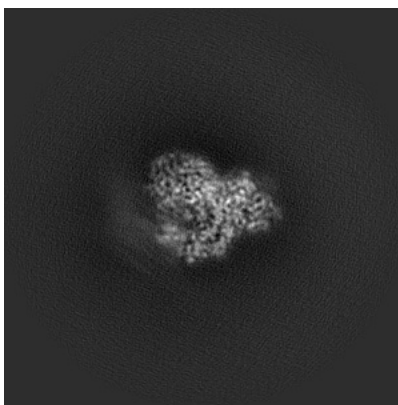


Z

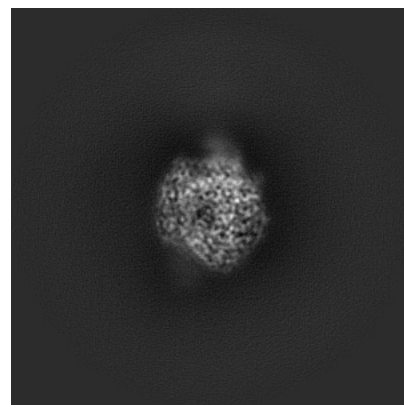
#### 6.1.2 Raw map



X



Y

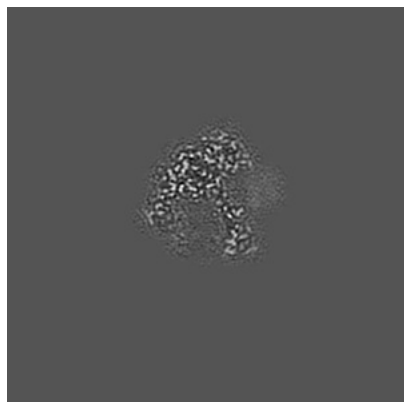


Z

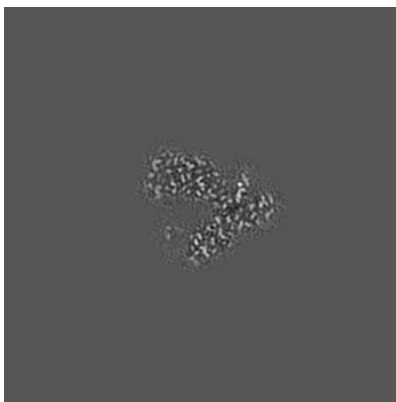
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

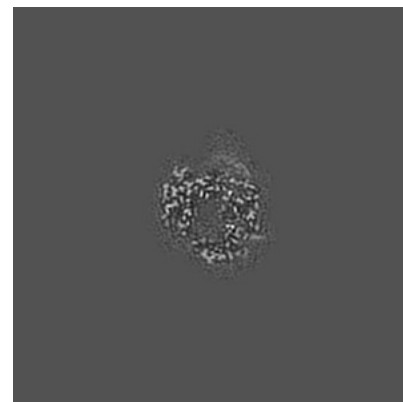
### 6.2.1 Primary map



X Index: 150

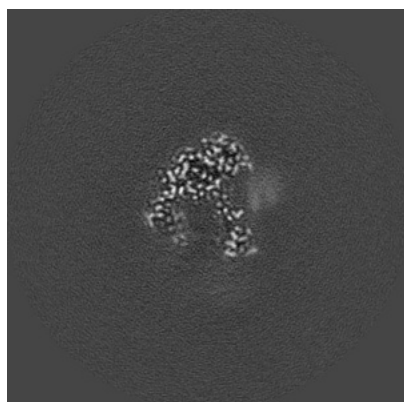


Y Index: 150

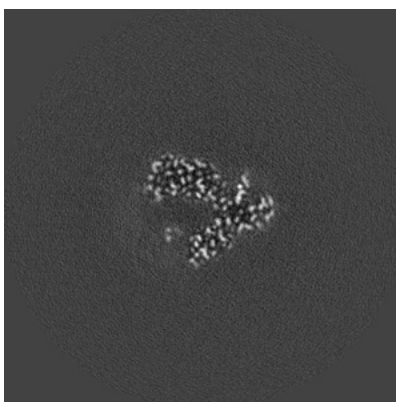


Z Index: 150

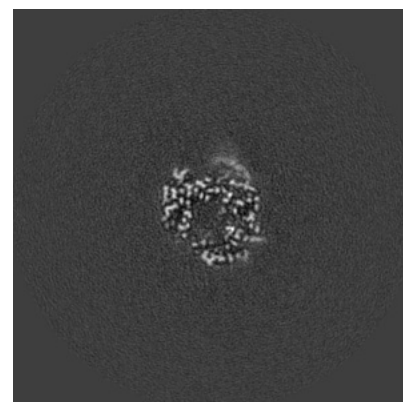
### 6.2.2 Raw map



X Index: 150



Y Index: 150

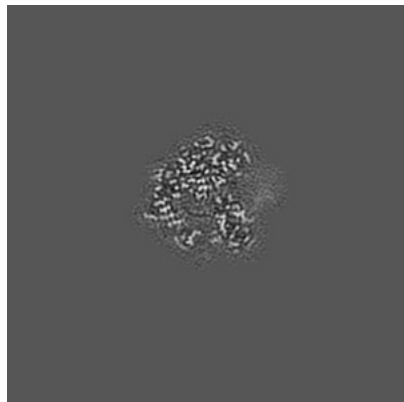


Z Index: 150

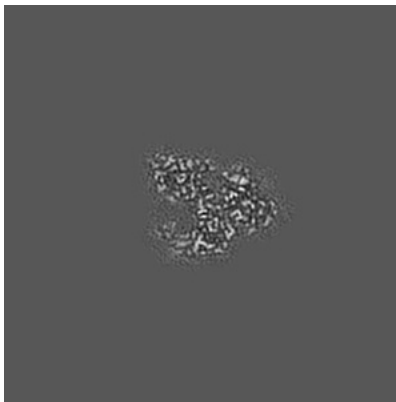
The images above show central slices of the map in three orthogonal directions.

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

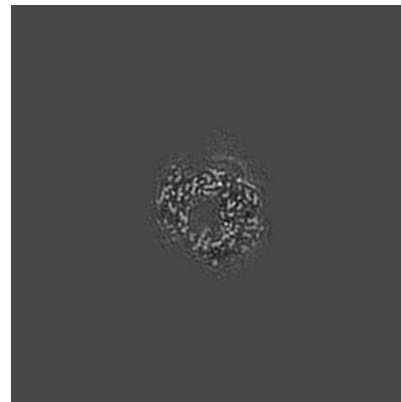
### 6.3.1 Primary map



X Index: 155

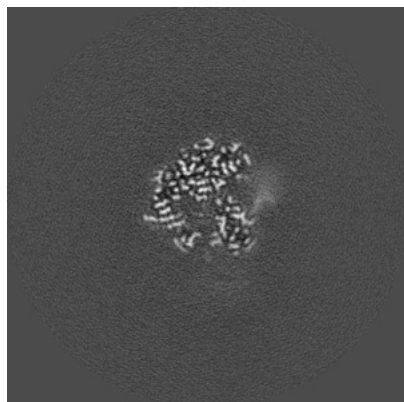


Y Index: 159

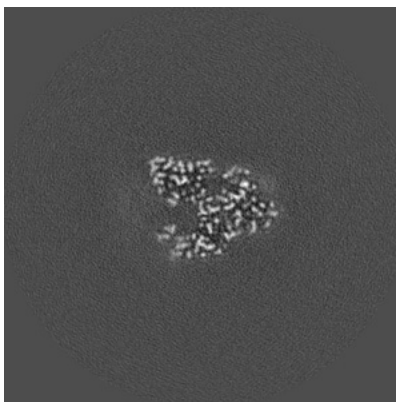


Z Index: 148

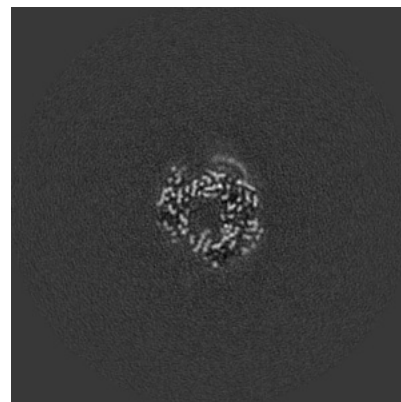
### 6.3.2 Raw map



X Index: 155



Y Index: 158

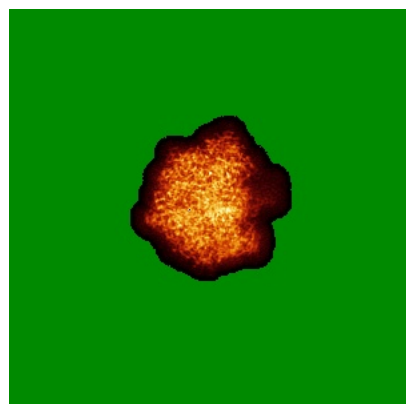


Z Index: 148

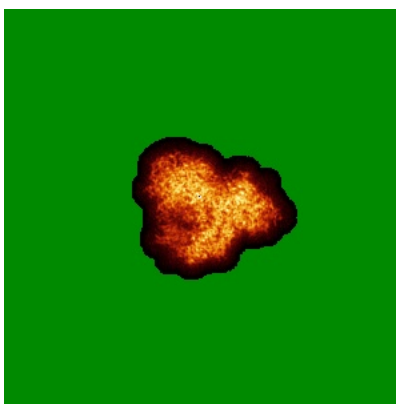
The images above show the largest variance slices of the map in three orthogonal directions.

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

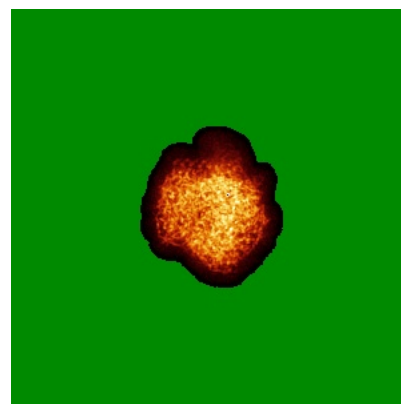
### 6.4.1 Primary map



X

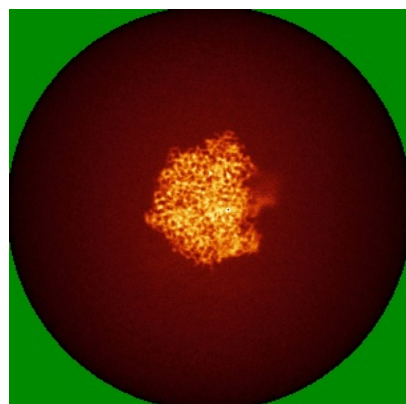


Y

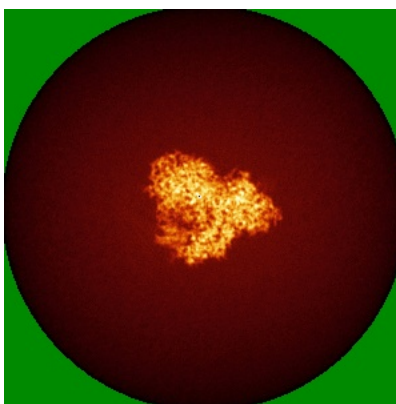


Z

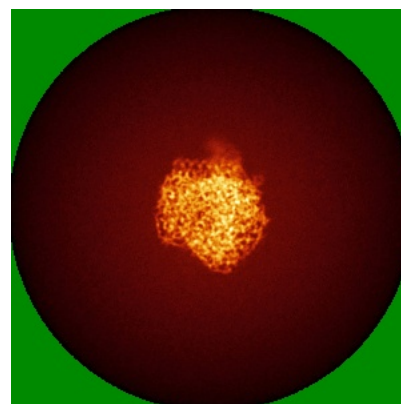
### 6.4.2 Raw map



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.

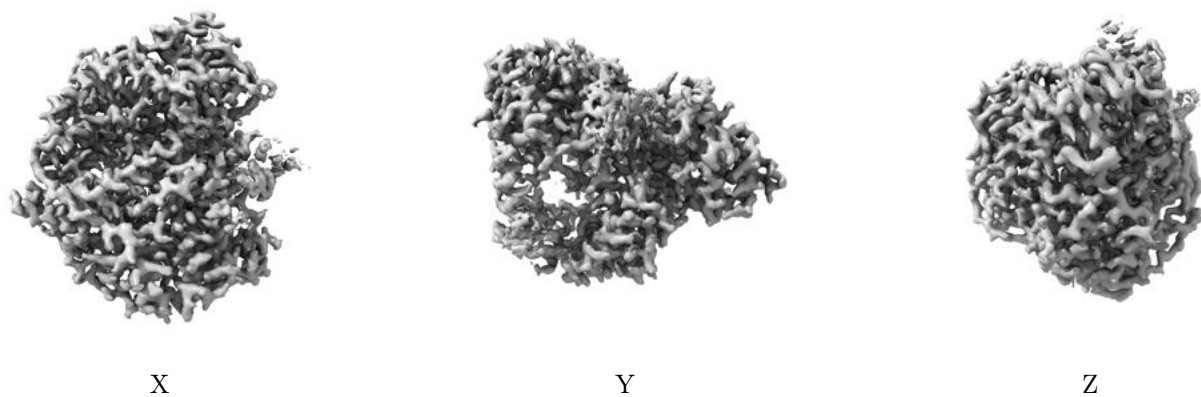
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.035. 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.

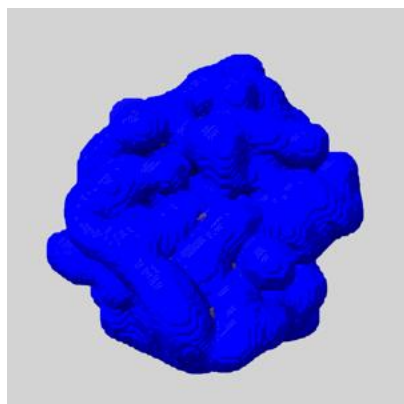
## 6.6 Mask visualisation [i](#)

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

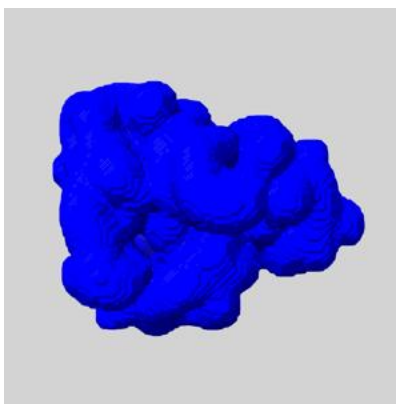
A mask typically either:

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

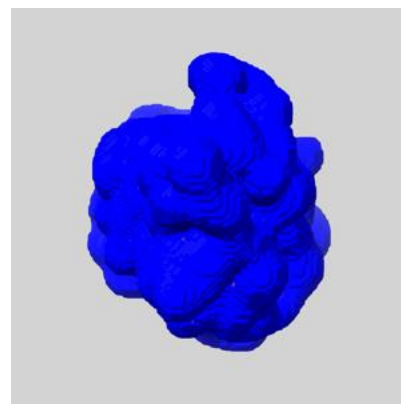
### 6.6.1 emd\_17858\_msk\_1.map [i](#)



X



Y

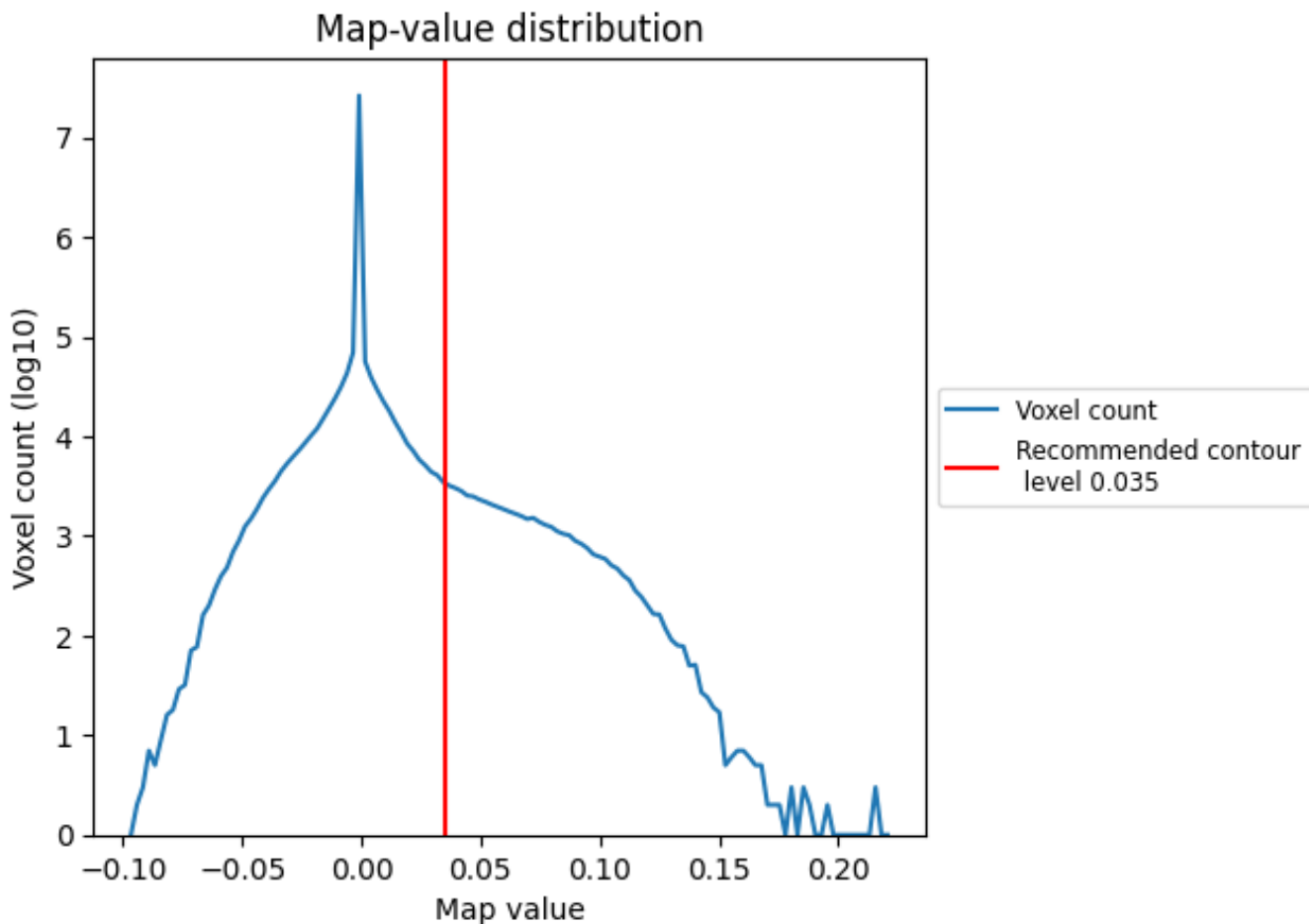


Z

## 7 Map analysis [i](#)

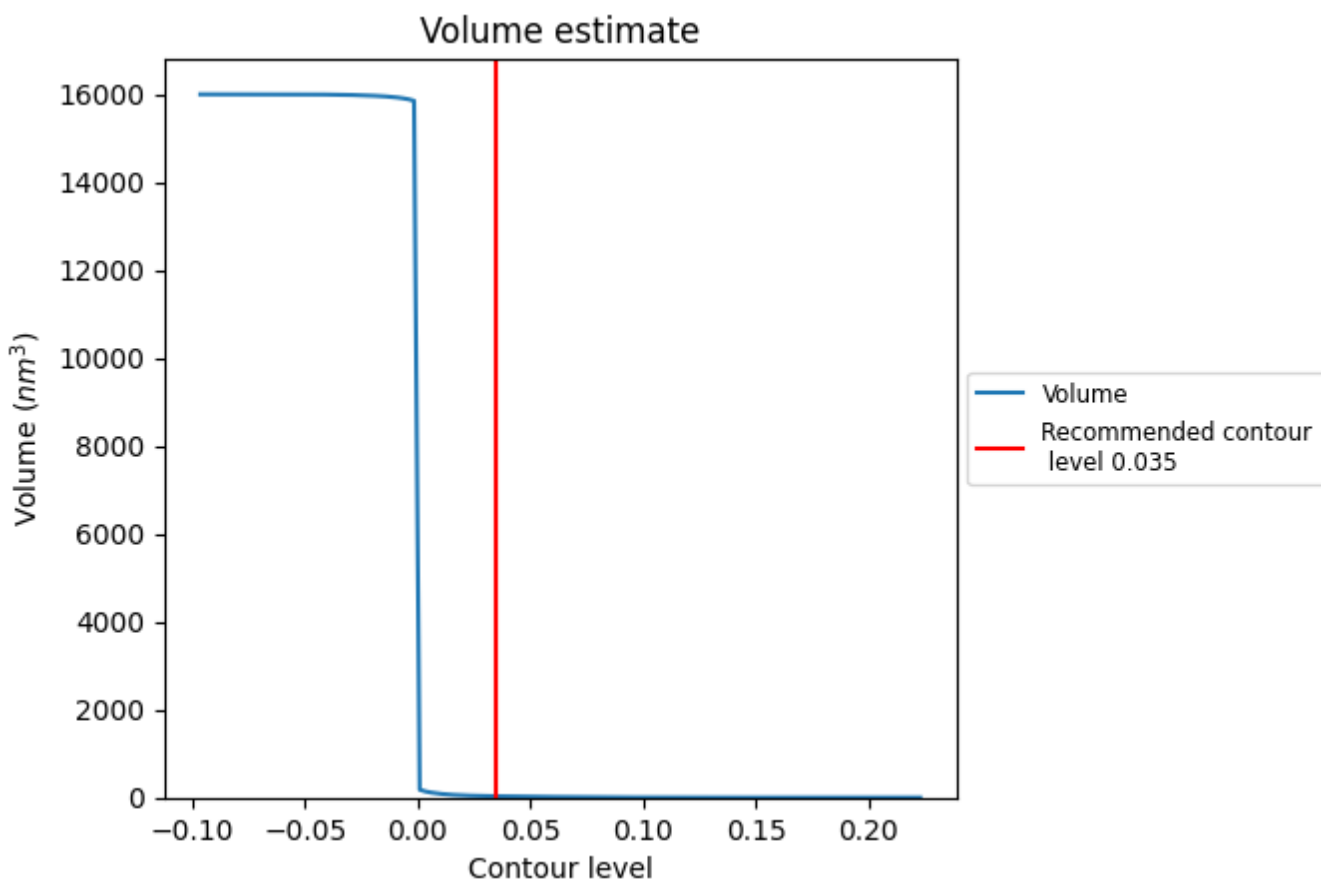
This section contains the results of statistical analysis of the map.

### 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.

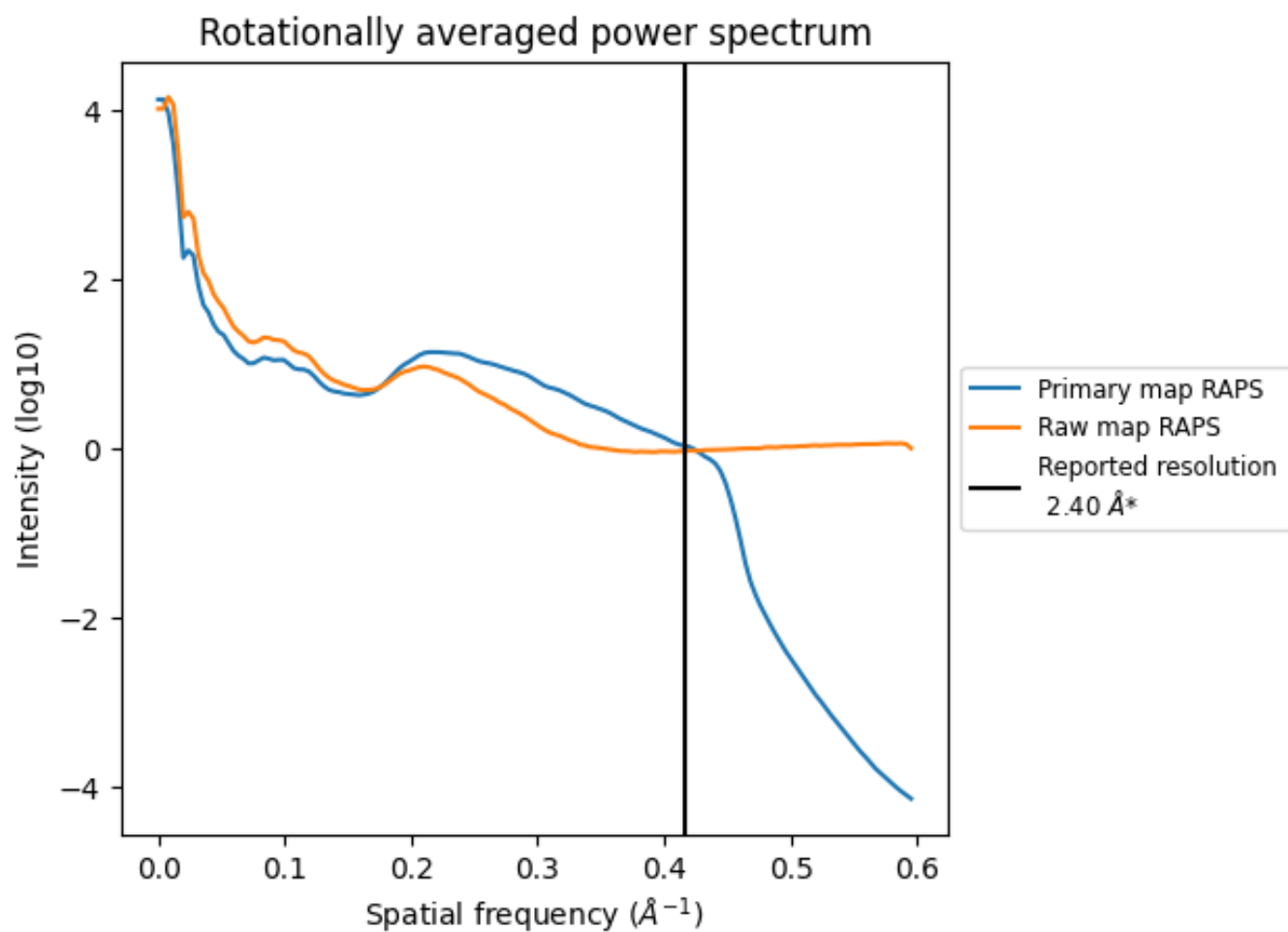
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 30 nm<sup>3</sup>; this corresponds to an approximate mass of 27 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](#)

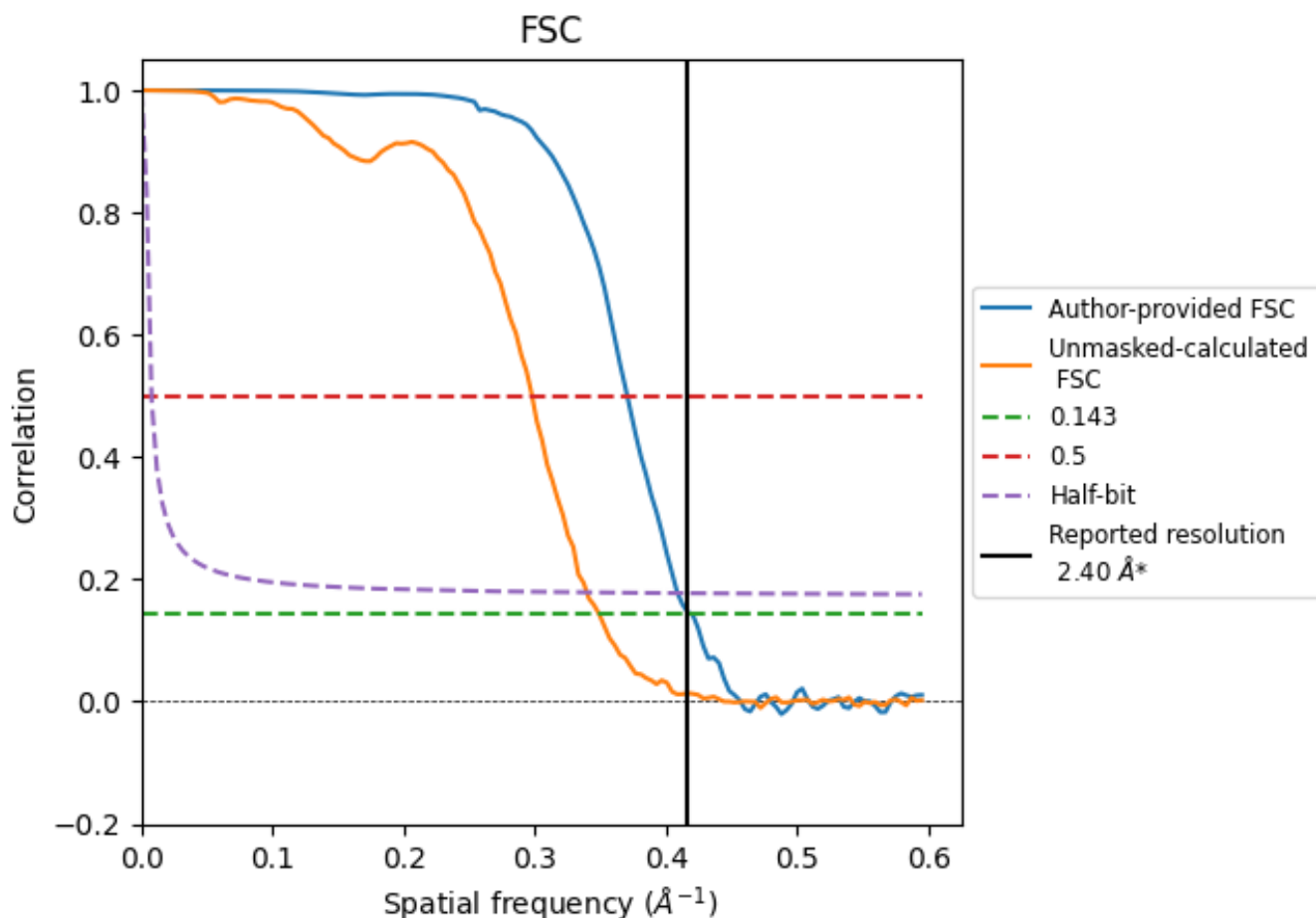


\*Reported resolution corresponds to spatial frequency of 0.417 Å<sup>-1</sup>

## 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  $0.417 \text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

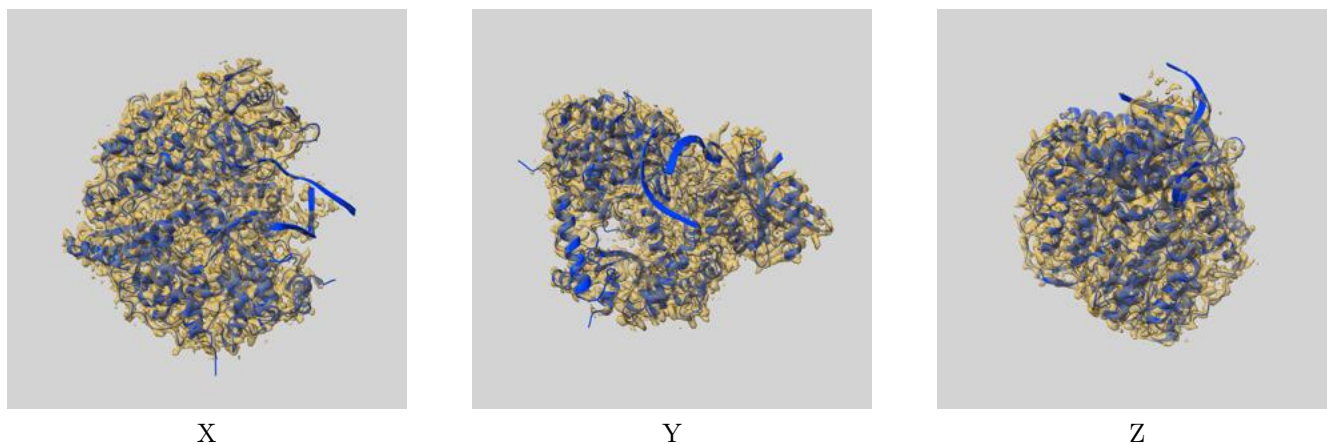
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.40	-	-
Author-provided FSC curve	2.38	2.70	2.45
Unmasked-calculated*	2.86	3.36	2.94

\*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 2.86 differs from the reported value 2.4 by more than 10 %

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

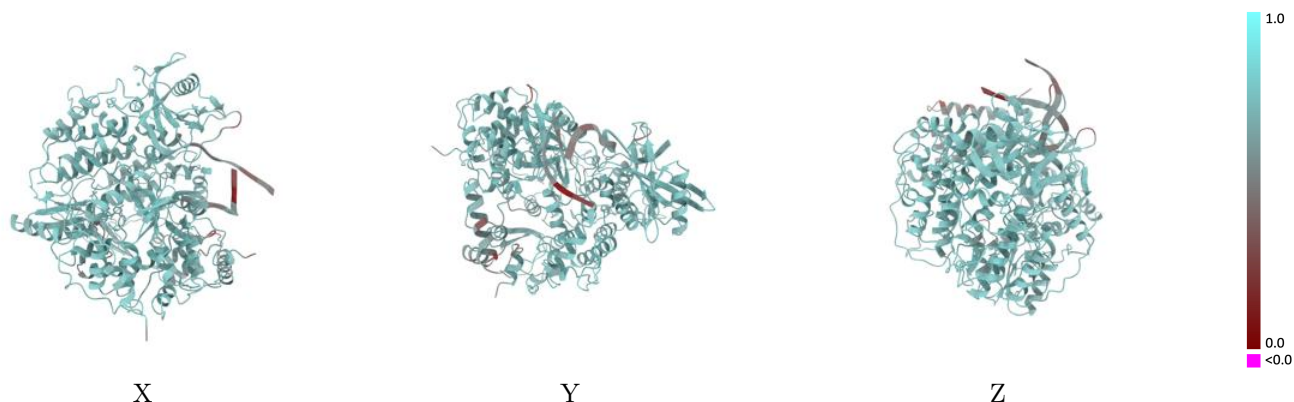
This section contains information regarding the fit between EMDB map EMD-17858 and PDB model 8PSO. Per-residue inclusion information can be found in section 3 on page 7.

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



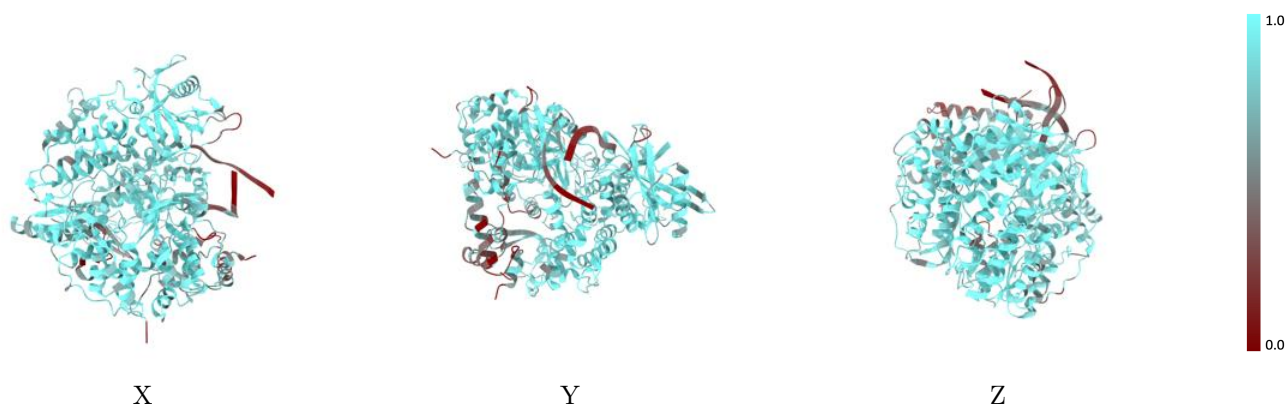
The images above show the 3D surface view of the map at the recommended contour level 0.035 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



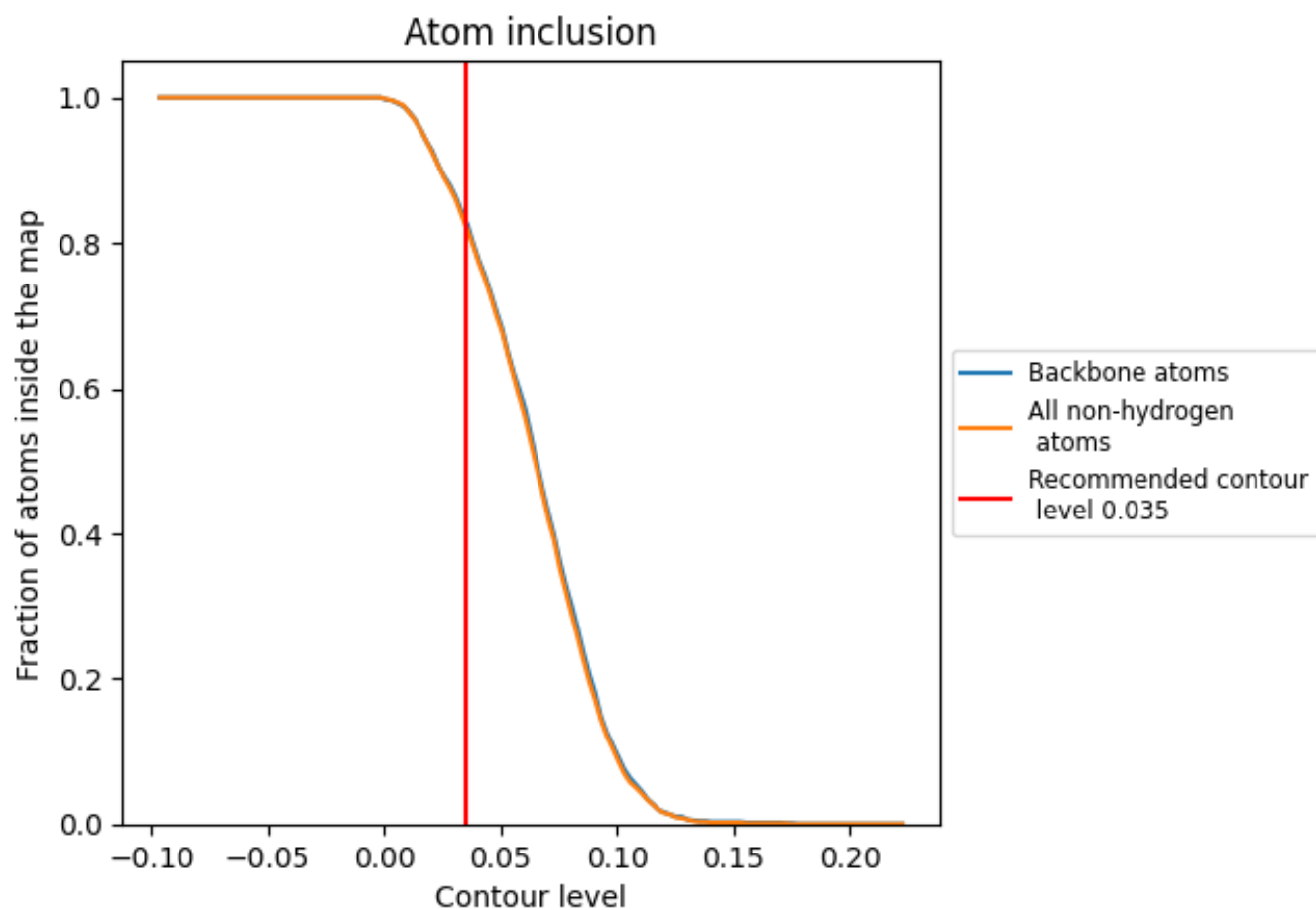
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.035).















## 9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 82% 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 (0.035) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8240	 0.6790
A	 0.8810	 0.6990
B	 0.8450	 0.6860
C	 0.7710	 0.6670
F	 0.9310	 0.6900
S	 0.6490	 0.5590
V	 0.6120	 0.5790

