

wwPDB EM Validation Summary Report (i)

Mar 7, 2024 - 01:12 pm GMT

PDB ID	:	8PT6
EMDB ID	:	EMD-17868
Title	:	Tilapia Lake Virus polymerase in vRNA initiation state (replicase conforma-
		tion)
Authors	:	Arragain, B.; Cusack, S.
Deposited on		
Resolution	:	3.03 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* A user guide is available at https://www.wwpdb.org/validation/2017/EMValidationReportHelp with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

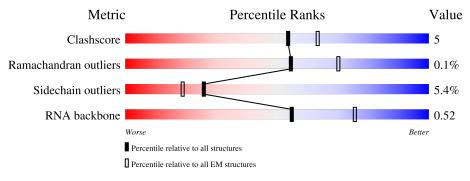
EMDB validation analysis	:	0.0.1.dev70
Mogul	:	1.8.4, CSD as 541 be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 3.03 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f EM} {f structures} \ (\#{f Entries})$
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 $\leq 40\%$). The numeric value is given above the bar.

Mol	Chain	Length		Quality of chain							
1	А	419	24%		909	26		10	%		
0			•								
2	В	519	22%		81%			17%	••		
3	С	478	2276	52%	6	12% •	35%				
4	S	40	8%	25%	5%	62%					
4	V	40	22%	8%) •	68%			_		
5	F	1				100%					



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 20448 atoms, of which 10063 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	А	417	Total 6652	C 2105	Н 3314	N 597	0 617	S 19	0	0

• Molecule 2 is a protein called Putative PB1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
2	В	515	Total 7971	C 2490	Н 3998	N 694	O 759	S 30	1	0

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

Mol	Chain	Residues	Atoms						AltConf	Trace
3	С	310	Total 4887	C 1563	Н 2440	N 432	0 438	S 14	0	0

There are 22 discrepancies between the modelled and reference sequences:

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

Continued on next page...



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

Continued from previous page...

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
4	V	13	Total	С	Η	Ν	Ο	Р	0	0
4	4 V	10	406	121	138	42	92	13	0	0
4	C	15	Total	С	Н	Ν	Ο	Р	0	0
4	4 5	19	486	145	161	60	105	15	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
5	F	1	Total 41	-	Н 12		-	Р 3	0	0

• Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	AltConf
6	А	1	Total Zn 1 1	0
6	С	2	Total Zn 2 2	0

• Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

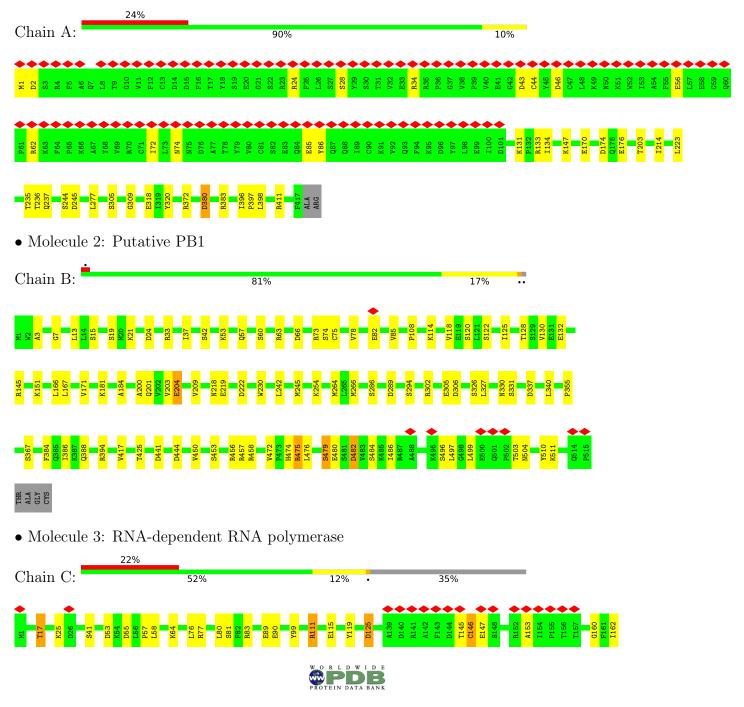
Mol	Chain	Residues	Atoms	AltConf
7	В	1	Total Mg 1 1	0
7	F	1	Total Mg 1 1	0



3 Residue-property plots (i)

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)



C163 C164 C164 C164 C164 C165 C168 C168 C168 C168 C170 C177 C177	K178 1179 6181 6181 6181 6182 6182 6182 6182 6182	V134 V135 E195 A198 A198 F200 F200 F200 A202 S203 S203 S205 F206 S203 S205 S205 S205 S205	R213 L214 A215 A215 L218 N219 V21 V221 V221 V223	E224 D225 P226 S227 V229 A230 A230 A231 A231 C233 C233
P234 C235 K2336 V236 V236 V238 V238 F244 F243 F243 F243 F243 F244 F244	A249 A245 F256 A257 F258 C258 C258 C258 C256 4 V264 V264 V264	H266 H267 P268 C270 C270 ALA ALA ALA ALA ALA ALA ALA ALA ALA AL	LEU PHE CLN GLN GLN ALA ALA VAL SEU ASN CLEU CLEU CLEU	GLN GLY SER ASN
VAL VAL SER CLY SER CLY SER HIS THR HIS ASP ASP ASP ARG ARG	ALA ALA TYR TYR SER ARG CLY ALA SER ALA SER CLU CLU CLU	LEU ALA ASIA ASIA ASIA THE THE PRO CLU CLY CTS CTS CTS CTS CTS CTS CTS CTS CTS CTS	ASP ARG LYS LYS LYS LEU AGN ALA ALA ALA SER SER SER	ARG
TYR THE THE THE THE THE THE CLU CLEU VAL VAL VAL ASN ASN ASN	VAL VAL VAL VAL ARC PHE TYR TYR TYR ASP X331 ASP X331 ASP X332 X333 X333 X333 X333 X333 X333 X33	L387 S388 S388 S390 R391 K391 K391 K391 K395 C396 D396 D398 L398 D398	7400 7400 7400 7403 7403 7405 7405 7405 7405 7409 7409	R410 K411 K413 K413 K413 M415 L414 L415 L417 L418 S419 S419
C420 SER SER SER SER C420 GLY THR THR THR THR THR LYS LYS LYS ARG CYAL	VALA VALA GLIM GLIM ASIP ARIG ARIG ARIG ASIP ARIG ASID ARIG ASID ARIG ARIG ARIG ARIG ARIG ARIG ARIG ARIG	VAL VAL ALA ALA ALA ALA ASP ASP ASP CLU CLN SER SER SER SER CLN CLN LEN	TYR PHE GLN GLN HIS HIS HIS HIS HIS HIS HIS	HIS
• Molecule 4: 5' v Chain V:	vRNA end - vRNA loc	pp (40-mer) 68%		
G 43 45 01 010 011 011 011 011 011 010 011 010	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	< < < < D D D 0 0		
	vRNA end - vRNA loc	op (40-mer)		
Chain S: 8%	25% 5% < 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	62% 7 0 7 0 7 1 7 1 7 1 7 1 7 1 7 1 7 1 7 1 7 1 7 1		
• Molecule 5: DN	A (5'-D(*(CTP))-3')			
Chain F:		100%		

There are no outlier residues recorded for this chain.



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	80175	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 $(e^-/\text{\AA}^2)$	40	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	0.049	Depositor
Minimum map value	-0.023	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.007	Depositor
Map size (Å)	251.99998, 251.99998, 251.99998	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles $(^{\circ})$	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, 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	
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.25	0/3413	0.50	0/4611
2	В	0.25	0/4050	0.50	0/5481
3	С	0.26	0/2508	0.50	0/3381
4	S	0.15	0/364	0.67	0/566
4	V	0.18	0/297	0.71	0/458
All	All	0.25	0/10632	0.51	0/14497

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	А	3338	3314	3313	24	0
2	В	3973	3998	3995	53	0
3	С	2447	2440	2440	42	0
4	S	325	161	161	11	0
4	V	268	138	139	1	0
5	F	29	12	12	0	0
6	А	1	0	0	0	0
6	С	2	0	0	0	0

Continued on next page...



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	В	1	0	0	0	0
7	F	1	0	0	0	0
All	All	10385	10063	10060	112	0

Continued from previous page...

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 5.

The worst 5 of 112 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)
3:C:266:HIS:CD2	3:C:269:VAL:HG13	1.87	1.08
3:C:266:HIS:HD2	3:C:269:VAL:HG13	0.97	1.08
2:B:218:ASN:ND2	2:B:219:GLU:OE2	2.09	0.86
1:A:56:GLU:N	1:A:56:GLU:OE2	2.11	0.82
1:A:74:ASN:OD1	2:B:475:ARG:NH1	2.24	0.71

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	Perce	ntiles
1	А	415/419~(99%)	401 (97%)	14(3%)	0	100	100
2	В	514/519~(99%)	505~(98%)	8 (2%)	1 (0%)	47	80
3	С	306/478~(64%)	282 (92%)	24 (8%)	0	100	100
All	All	1235/1416~(87%)	1188 (96%)	46 (4%)	1 (0%)	54	84

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	В	254	LYS



5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Perce	ntiles
1	А	364/365~(100%)	355~(98%)	9~(2%)	47	77
2	В	445/447~(100%)	413 (93%)	32 (7%)	14	43
3	С	262/405~(65%)	245~(94%)	17 (6%)	17	47
All	All	1071/1217~(88%)	1013 (95%)	58 (5%)	26	55

5 of 58 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
2	В	337	ASP
3	С	381	LYS
2	В	479	SER
3	С	267	ARG
3	С	111	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	371	HIS
3	С	266	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	S	15/40~(37%)	2(13%)	1 (6%)
4	V	12/40~(30%)	4 (33%)	0
All	All	27/80~(33%)	6 (22%)	1 (3%)

5 of 6 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	V	5	А
4	V	7	С

Continued on next page...



Continued from previous page...

Mol	Chain	Res	Type
4	V	10	U
4	V	11	С
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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 5 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.



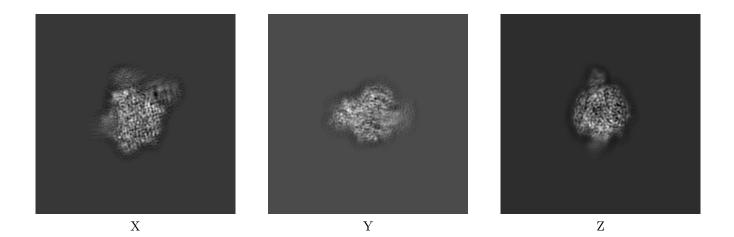
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-17868. These allow visual inspection of the internal detail of the map and identification of artifacts.

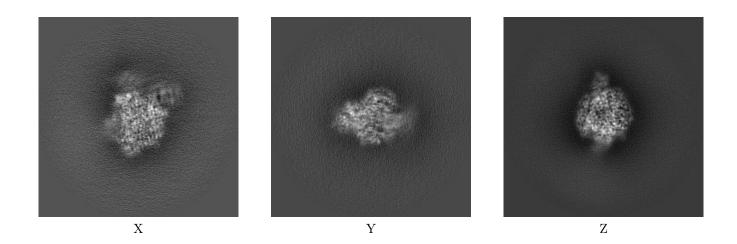
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

6.1.1 Primary map



6.1.2 Raw map



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

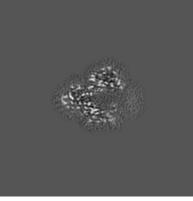


6.2 Central slices (i)

6.2.1 Primary map



X Index: 150

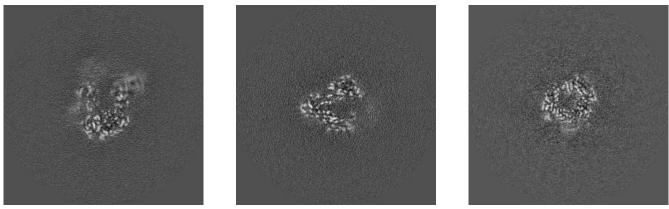


Y Index: 150



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.

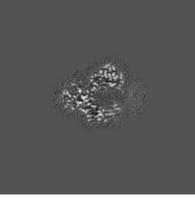


6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 153

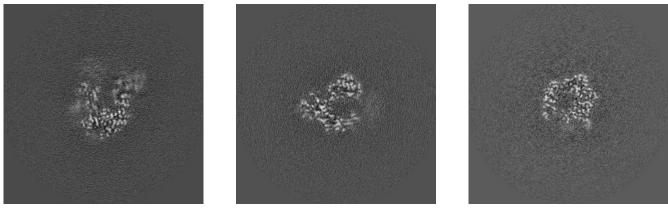


Y Index: 147



Z Index: 156

6.3.2 Raw map



X Index: 153

Y Index: 147

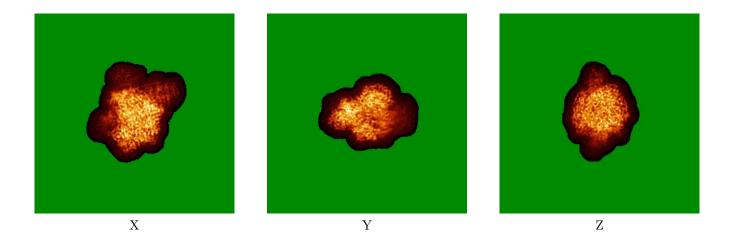


The images above show the largest variance slices of the map in three orthogonal directions.

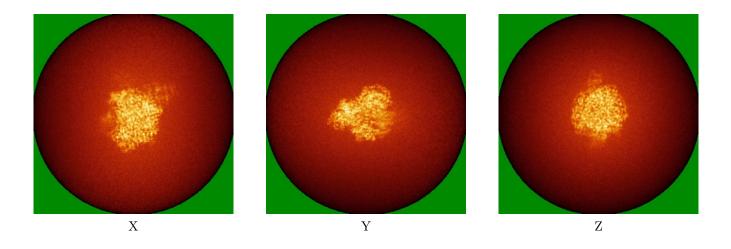


6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



6.4.2 Raw map



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



6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.007. 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.



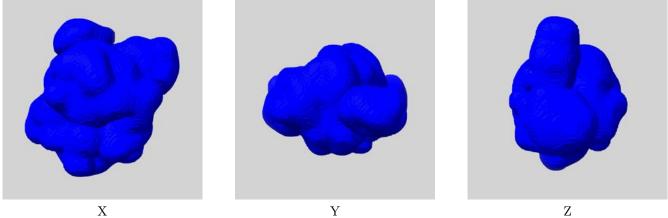
Mask visualisation (i) 6.6

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

$emd_{17868}msk_{1.map}$ (i) 6.6.1

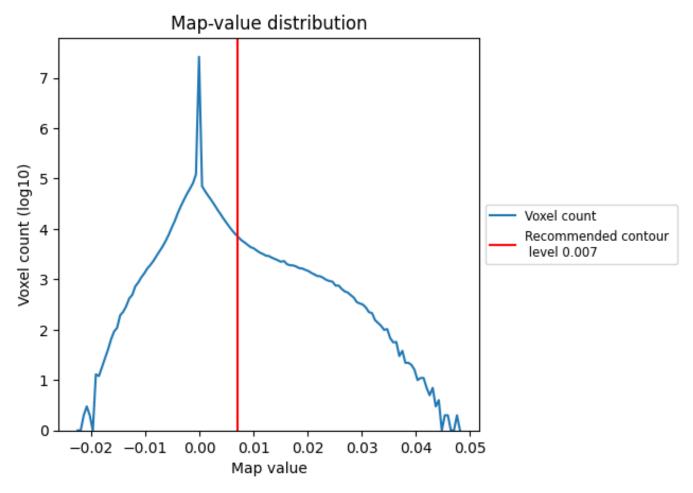




7 Map analysis (i)

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

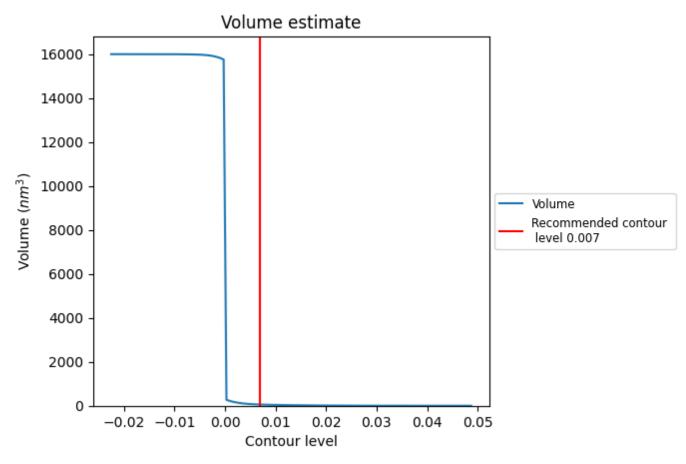
7.1 Map-value distribution (i)



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)

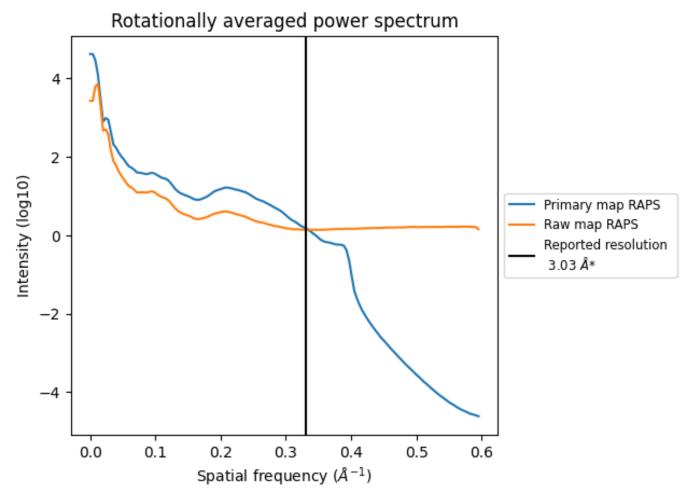


The volume at the recommended contour level is 54 nm^3 ; this corresponds to an approximate mass of 49 kDa.

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)



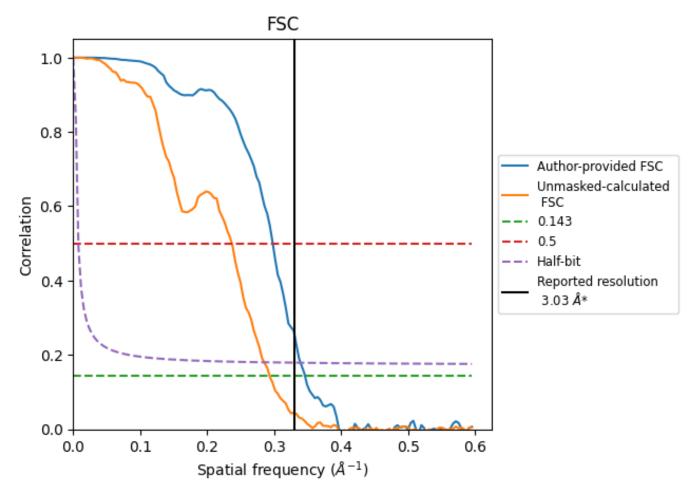
*Reported resolution corresponds to spatial frequency of 0.330 ${\rm \AA^{-1}}$



8 Fourier-Shell correlation (i)

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

8.1 FSC (i)



*Reported resolution corresponds to spatial frequency of 0.330 ${\rm \AA^{-1}}$



8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off		
Resolution estimate (A)	0.143	0.5	Half-bit
Reported by author	3.03	-	-
Author-provided FSC curve	2.89	3.36	2.95
Unmasked-calculated*	3.41	4.22	3.51

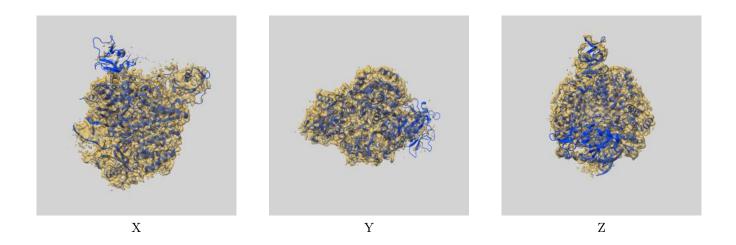
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.41 differs from the reported value 3.03 by more than 10 %



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-17868 and PDB model 8PT6. Per-residue inclusion information can be found in section 3 on page 5.

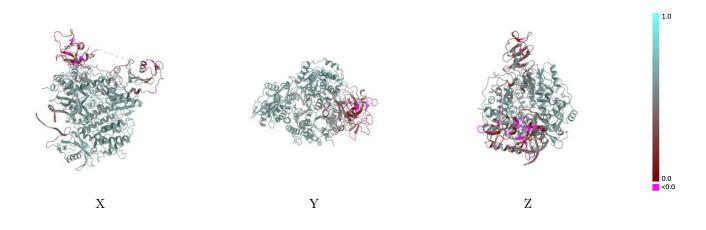
9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.007 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.

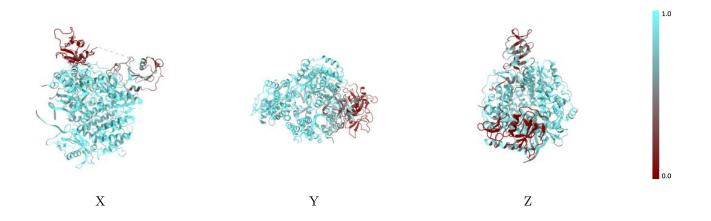


9.2 Q-score mapped to coordinate model (i)



The images above show the model with each residue coloured according 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.

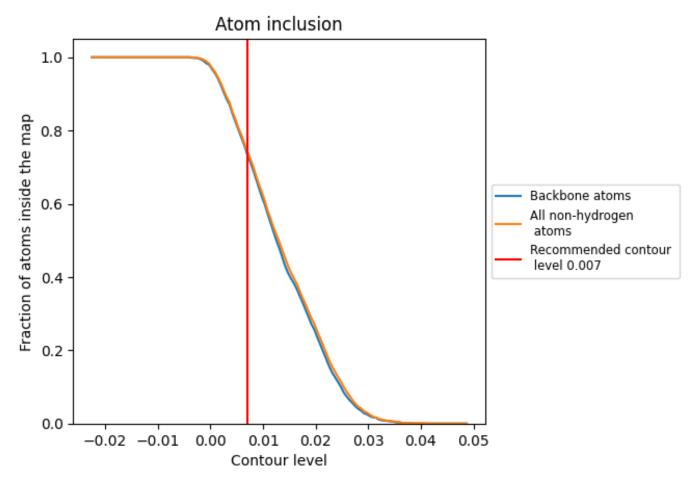
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.007).



9.4 Atom inclusion (i)



At the recommended contour level, 74% of all backbone atoms, 74% of all non-hydrogen atoms, are inside the map.



1.0

0.0 <0.0

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.007) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7430	0.5130
А	0.6930	0.4920
В	0.8910	0.5900
С	0.5740	0.4230
F	0.9000	0.5710
S	0.8250	0.5030
V	0.8280	0.4760

