



wwPDB EM Validation Summary Report ⓘ

Jun 29, 2026 – 04:04 pm BST

PDB ID : 9TIO / pdb_00009tio
EMDB ID : EMD-55963
Title : Phage 812 baseplate in the pre-contraction state - upper arm (segment B)
Authors : Binovsky, J.; Plevka, P.
Deposited on : 2025-12-05
Resolution : 3.90 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 ⓘ 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.dev133
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

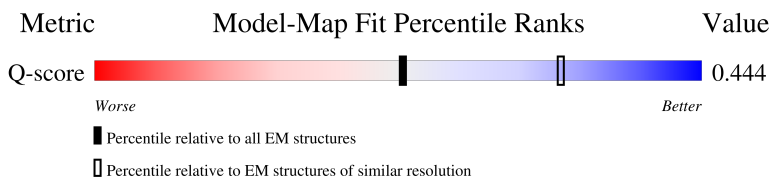
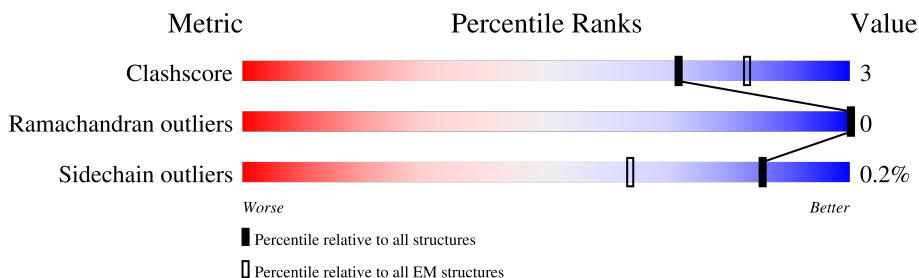
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 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	-
Q-score	-	25397	8855 (3.40 - 4.40)

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	1152	
1	B	1152	
1	D	1152	
2	C	173	

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Mol	Chain	Length	Quality of chain
2	E	173	89%10%
2	F	173	87%7%6%
2	K	173	14%76%11%12%
3	G	1019	22%75%
4	H	640	21%77%
4	I	640	21%75%
4	J	640	22%76%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 15828 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 protein called ORF65.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	342	Total	C	N	O	S	0	0
			2695	1692	441	554	8		
1	B	165	Total	C	N	O	S	0	0
			1306	823	211	269	3		
1	D	166	Total	C	N	O	S	0	0
			1314	829	212	270	3		

- Molecule 2 is a protein called ORF64.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	111	Total	C	N	O	S	0	0
			889	578	145	165	1		
2	E	172	Total	C	N	O	S	0	0
			1349	858	221	269	1		
2	F	163	Total	C	N	O	S	0	0
			1277	817	208	251	1		
2	K	152	Total	C	N	O	S	0	0
			1199	769	196	233	1		

- Molecule 3 is a protein called ORF63.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	257	Total	C	N	O	S	0	0
			2138	1364	339	428	7		

- Molecule 4 is a protein called CBM-cenC domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	145	Total	C	N	O	S	0	0
			1193	760	192	239	2		
4	I	157	Total	C	N	O	S	0	0
			1259	798	205	254	2		

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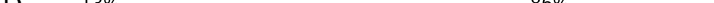
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	J	151	1209	768	198	241	2	0	0

- Molecule 1: ORF65



LYS SER THR THR ARG ALA ASN GLU PHE THR ARG TYR VAL ILE ASP CLU LYS VAL LYS SER SER GLY THR THR ASN LYS LEU LEU VAL ARG ASP LEU LEU LEU LEU ARG PRO VAL ARG ARG ARG MET VAL THR THR ARG ASP

- Molecule 1: ORF65

Chain B:  13% 86%

[illegible]

ARG	THR	TYR	ASP	GLU	SER	SER	TYR	LYS	VAL	ASN	GLY	PHE	GLU	LEU	PHE	SER	GLU	GLY	ASN	ALA	GLU	ASP	ASP	HIS	VAL	SER	VAL	VAL	VAL	ASP	ALA	GLY	LYS	LYS	PHE	LYS	VAL	VAL	ASP	LYS	PRO	VAL	SER	THR	ARG	ILE	SER	VAL	PRO	LYS	TYR	SER	ASP	LEU
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TLE	VAL	LVS	VAL	TRP	THR	GLU	THR	SER	PRO	GLY	VAL	THR	THR	LVS	GLU	LYS	TYR	LYS	GLN	ASP	PHE	ARG	LEU	THR	THR	ASP	GLY	GLN	THR	ILE	ASP	TRP	SER	PRO	GLN	GLY	GLN	GLU	GLU	PRO	PRO	SER	GLY	GLY	THR	SER	TYR	TYR	VAL	SER	LYS	TYR	LYS	MET	ALA	GLY	GLY
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| TRP | ASP | THR | GLU | ASN | VAL | THR | THR | GLU | GLN | LYS | LYS | LYS | VAL | THR | MET | LYS | ARG | PHE | TRP | ARG | ARG | HIS | ASN | GLU | GLY | TYR | TYR | GLY | GLU | THR | GLU | HIS | TYR | LEU | TYR | SER | SER | ASN | ASN | LEU | GLN | LEU | ASP | ALA | GLY | GLN | LYS | TRP | LYS | GLY | GLU | THR | TYR | ALA | TYR | ASP | ARG | GLU | HIS |

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

GLY	SER	GLU	ASP	GLY	THR	ILE	MET	THR	ASP	ALA	LYS	GLY	ALA	THR	LYS	GLY	PHE	THR	ILE	PRO	ALA	GLY	ILE	ARG	CYS	GLY	ASN	ARG	GLU	VAL	LEU	THR	LYS	ASN	ALA	ASN	SER	THR	SER	ALA	THR	THR	Tyr	THR	THR	THR	ALA	GLN	GLY	ARG	LYS	LYS	THR	ALA	ALA	GLN	ASP	ILE	ILE	ILE	ILE	ILE
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THR	VAL	MET	MET	ASN	ALA	ASP	ILE	VAL	SER	ASN	ASN	ALA	SER	ALA	GLU	THR	ARG	VAL	TYR	PHE	ASP	PRO	MET	MET	MET	GLY	GLY	LYS	GLU	TYR	THR	ILE	VAL	ILE	ILE	THR	GLU	ASN	SER	ASP	TYR	THR	THR	MET	TRP	MET	VAL	GLY	THR	ARG	LYS	THR	LYS	PRO	LYS	ILE	ASP	LYS	PRO
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ASN	GLU	VAL	VAL	ILE	SER	GLY	ASN	ASN	PRO	TYR	LEU	GLN	GLY	VAL	LEU	PHE	SER	SER	SER	ASN	ALA	ALA	SER	THR	TRP	TRP	PRO	PRO	HIS	GLN	ASN	SER	SER	ASP	LEU	LYS	PHE	GLY	ILE	TYR	THR	SER	LYS	PHE	ASN	GLU	THR	ALA	THR	ILE	GLU	PHE	GLU	ILE	ILE	ASP	LYS	VAL	SER	ALA	ASP	ARG	ILE	VAL
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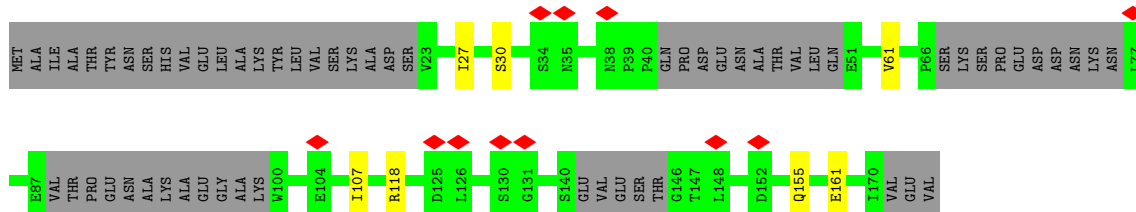
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- Molecule 1: ORF65

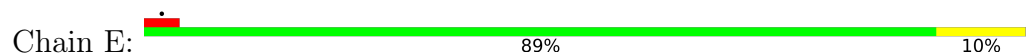


SER	ALA	ASP	ARG	ILE	VAL	LEU	MET	SER	THR	TYR	LEU	THR	PRO	GLU	ARG	THR	GLY	CYS	THR	TRP	GLU	MET	LYS	LEU	ILE	LEU	ASP	ASP	MET	ALA	SER	SER	THR	THR	PHE	ASP	GLN	LEU	LYS	TRP	GLU	GLU	PRO	ILE	GLY	ASN	GLN	VAL	LEU	ASP	ASP	VAL	GLY	LEU	ALA	ARG	GLN	VAL	LYS
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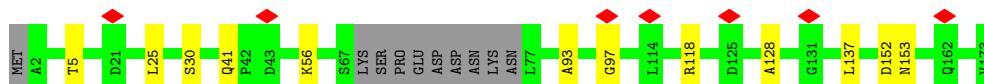
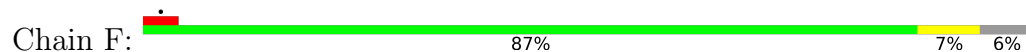
- Molecule 2: ORF64



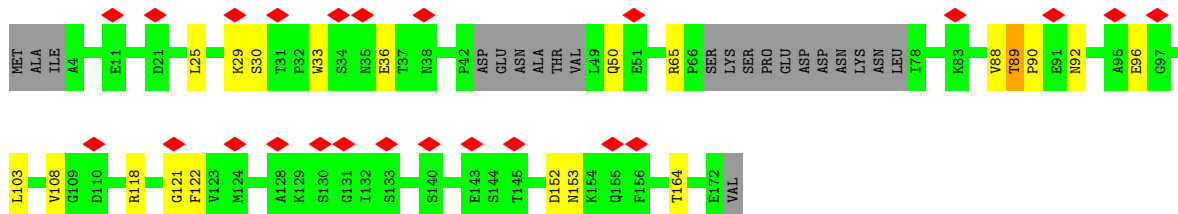
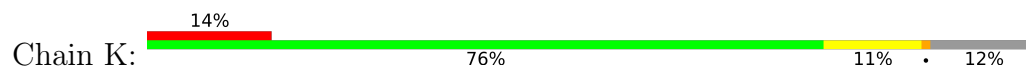
- Molecule 2: ORF64



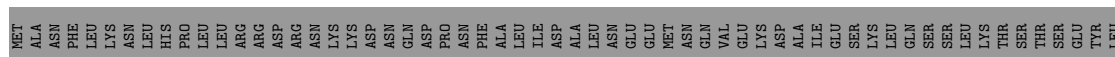
- Molecule 2: ORF64



- Molecule 2: ORF64

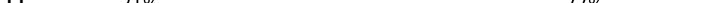


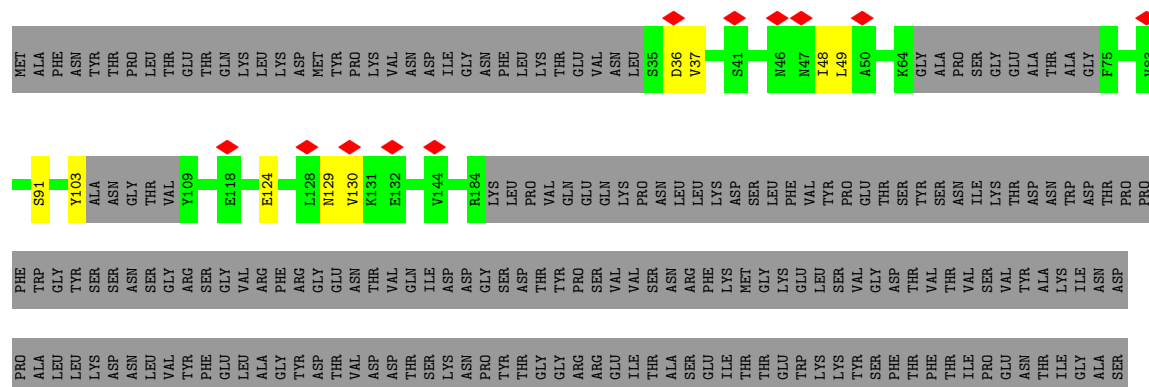
- Molecule 3: ORF63





- Molecule 4: CBM-cenC domain-containing protein

Chain H:  21% 77%



[illegible]

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	19727	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40.8	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.221	Depositor
Minimum map value	-0.111	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.04	Depositor
Map size ($\text{\AA}$)	300.096, 300.096, 300.096	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.6672001, 1.6672001, 1.6672001	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.11	0/2748	0.30	0/3725
1	B	0.10	0/1330	0.31	0/1800
1	D	0.12	0/1338	0.33	0/1811
2	C	0.11	0/907	0.29	0/1225
2	E	0.11	0/1377	0.32	0/1872
2	F	0.11	0/1303	0.28	0/1771
2	K	0.12	0/1224	0.30	0/1660
3	G	0.10	0/2185	0.29	0/2961
4	H	0.11	0/1219	0.30	0/1651
4	I	0.11	0/1288	0.27	0/1749
4	J	0.11	0/1236	0.32	0/1678
All	All	0.11	0/16155	0.30	0/21903

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	2695	0	2590	17	0
1	B	1306	0	1267	10	0
1	D	1314	0	1278	14	0
2	C	889	0	893	5	0
2	E	1349	0	1339	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1277	0	1274	7	0
2	K	1199	0	1195	12	0
3	G	2138	0	2032	18	0
4	H	1193	0	1144	5	0
4	I	1259	0	1207	13	0
4	J	1209	0	1164	5	0
All	All	15828	0	15383	105	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 3.

The worst 5 of 105 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)
1:B:83:ILE:O	1:B:88:ARG:NH1	2.18	0.77
1:B:69:SER:OG	1:B:71:ASP:OD1	2.05	0.75
1:A:88:ARG:NH2	1:A:143:GLU:OE1	2.20	0.74
2:F:30:SER:O	2:F:118:ARG:NH1	2.25	0.70
4:I:85:TYR:OH	4:I:105:ASN:OD1	2.05	0.69

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	336/1152 (29%)	333 (99%)	3 (1%)	0	100	100
1	B	163/1152 (14%)	162 (99%)	1 (1%)	0	100	100
1	D	164/1152 (14%)	162 (99%)	2 (1%)	0	100	100
2	C	101/173 (58%)	98 (97%)	3 (3%)	0	100	100
2	E	170/173 (98%)	164 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	159/173 (92%)	157 (99%)	2 (1%)	0	100	100
2	K	146/173 (84%)	142 (97%)	4 (3%)	0	100	100
3	G	251/1019 (25%)	247 (98%)	4 (2%)	0	100	100
4	H	137/640 (21%)	133 (97%)	4 (3%)	0	100	100
4	I	153/640 (24%)	153 (100%)	0	0	100	100
4	J	145/640 (23%)	141 (97%)	4 (3%)	0	100	100
All	All	1925/7087 (27%)	1892 (98%)	33 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	302/1010 (30%)	301 (100%)	1 (0%)	86	85
1	B	147/1010 (15%)	147 (100%)	0	100	100
1	D	148/1010 (15%)	147 (99%)	1 (1%)	76	78
2	C	100/153 (65%)	100 (100%)	0	100	100
2	E	152/153 (99%)	152 (100%)	0	100	100
2	F	143/153 (94%)	143 (100%)	0	100	100
2	K	134/153 (88%)	133 (99%)	1 (1%)	76	78
3	G	241/928 (26%)	241 (100%)	0	100	100
4	H	138/577 (24%)	137 (99%)	1 (1%)	76	78
4	I	142/577 (25%)	142 (100%)	0	100	100
4	J	136/577 (24%)	136 (100%)	0	100	100
All	All	1783/6301 (28%)	1779 (100%)	4 (0%)	85	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	936	PHE
1	D	92	ASN
4	H	103	TYR
2	K	89	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	157	GLN
3	G	201	GLN
4	J	397	ASN
4	J	42	GLN
4	J	172	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

There are no ligands in this entry.

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.

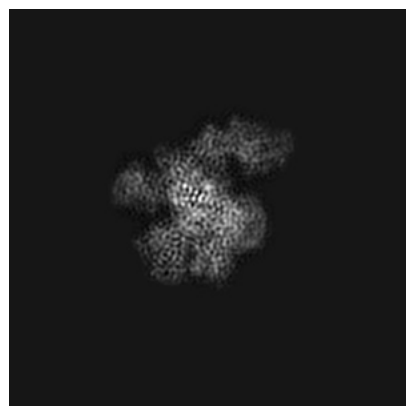
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55963. These allow visual inspection of the internal detail of the map and identification of artifacts.

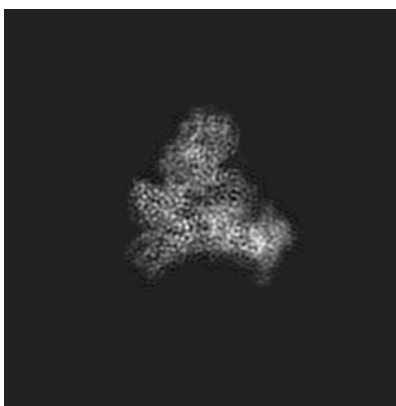
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

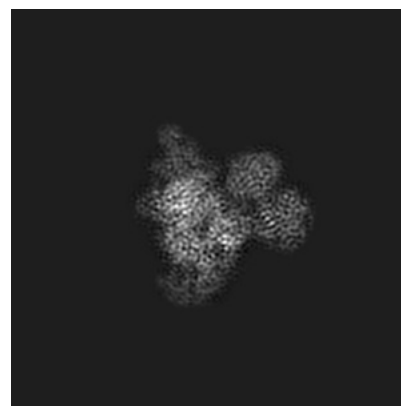
6.1.1 Primary map



X

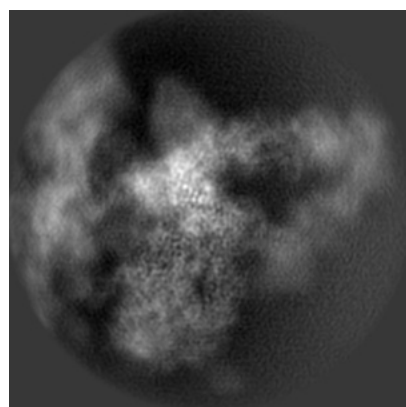


Y

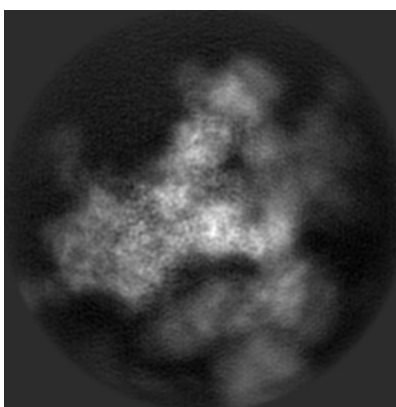


Z

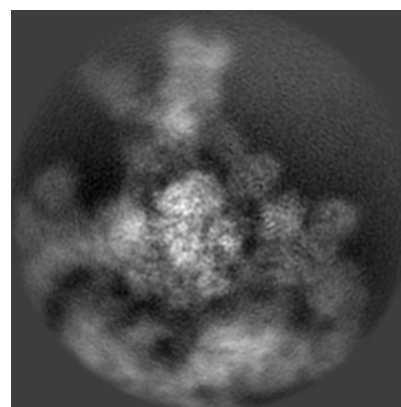
6.1.2 Raw map



X



Y

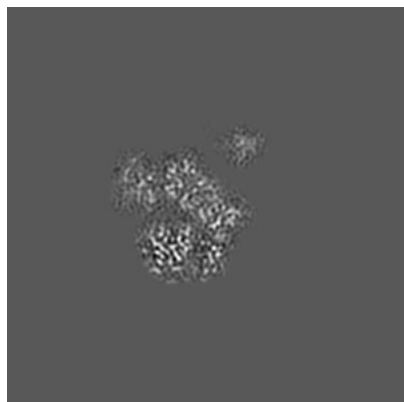


Z

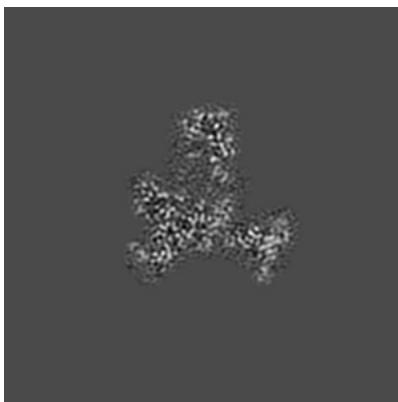
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

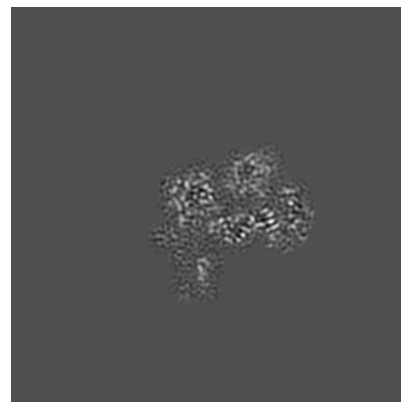
6.2.1 Primary map



X Index: 90

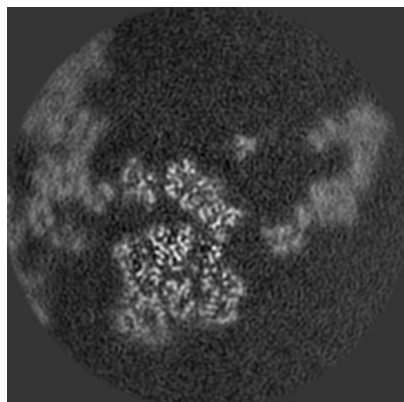


Y Index: 90

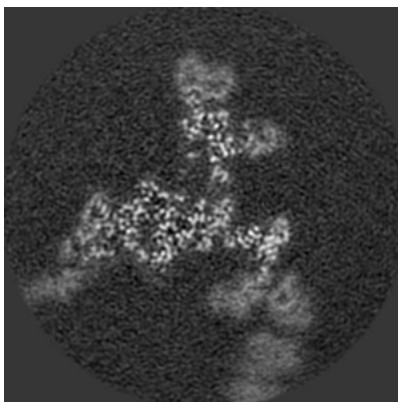


Z Index: 90

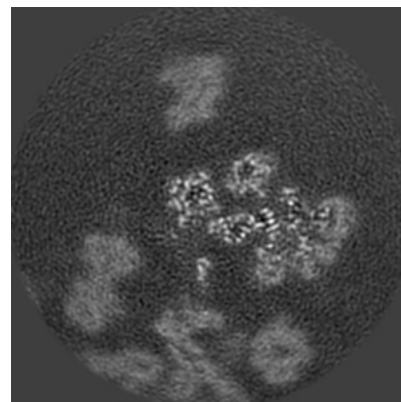
6.2.2 Raw map



X Index: 90



Y Index: 90

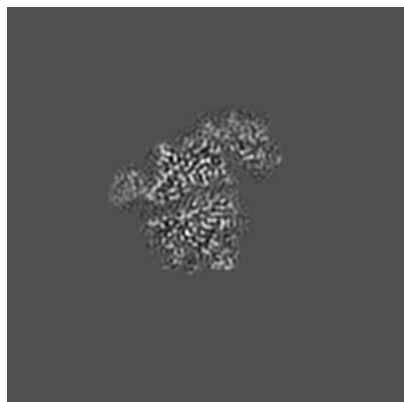


Z Index: 90

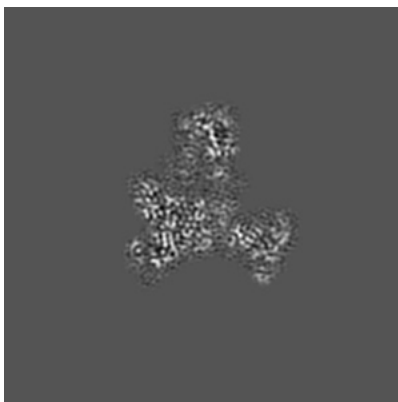
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

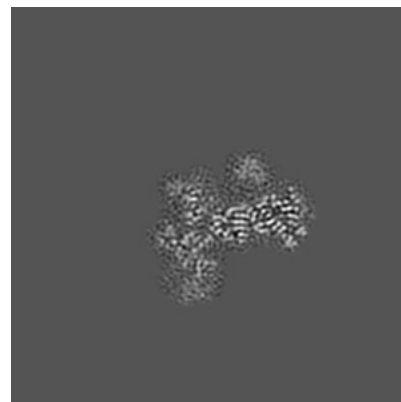
6.3.1 Primary map



X Index: 78

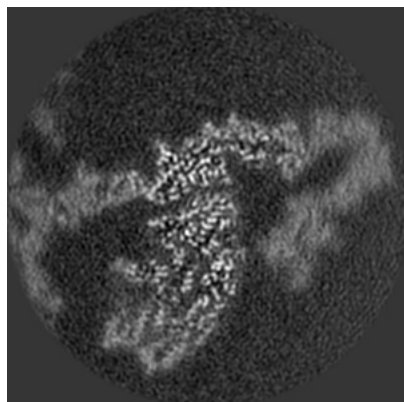


Y Index: 91

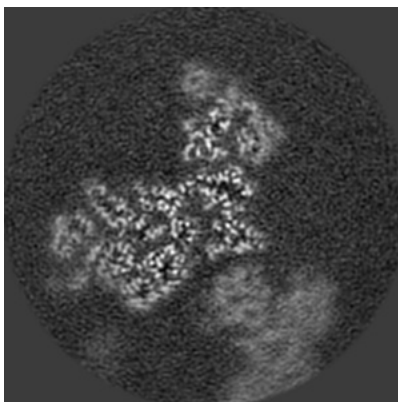


Z Index: 94

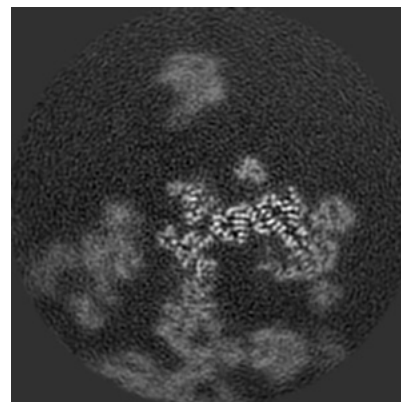
6.3.2 Raw map



X Index: 78



Y Index: 78

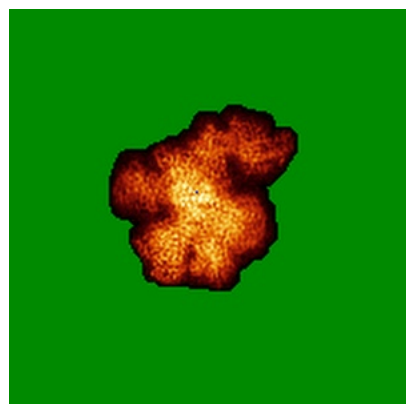


Z Index: 94

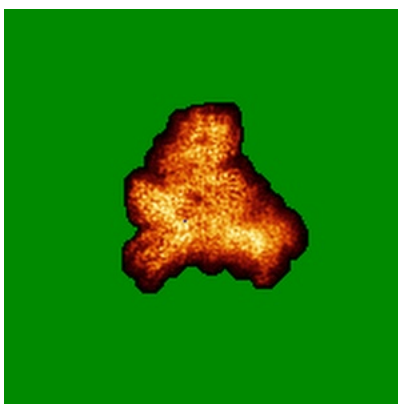
The images above show the largest variance slices of the map in three orthogonal directions.

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

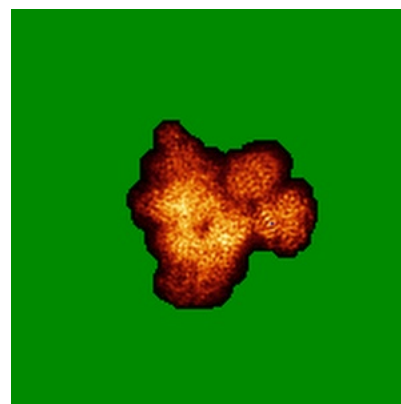
6.4.1 Primary map



X

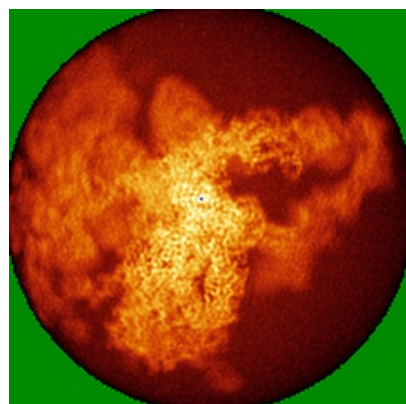


Y

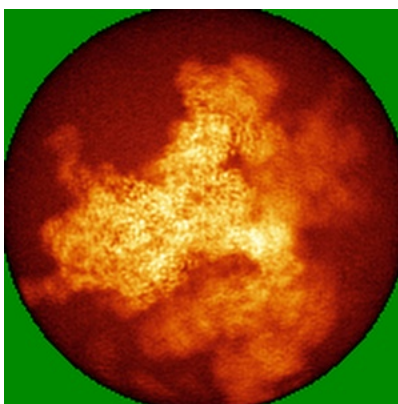


Z

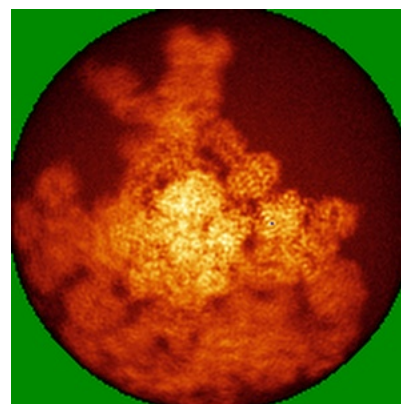
6.4.2 Raw map



X



Y

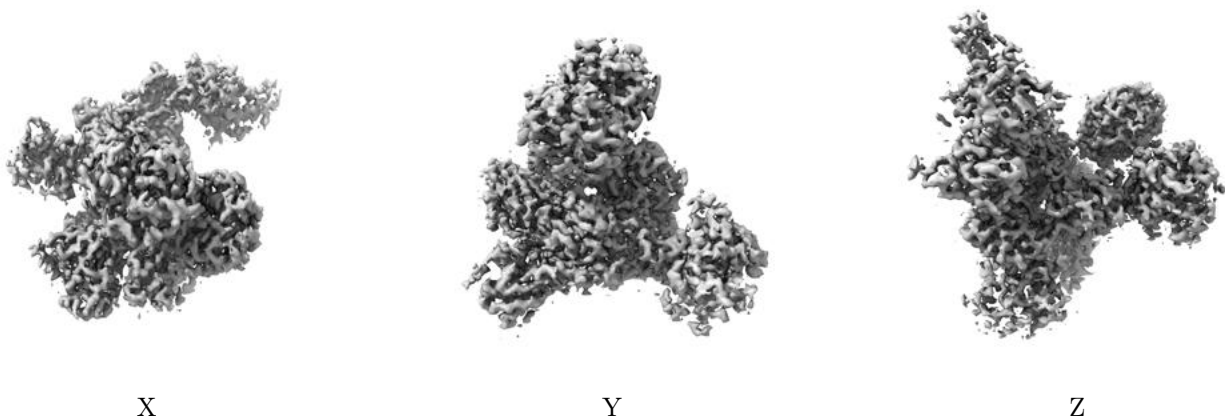


Z

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 0.04. 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.

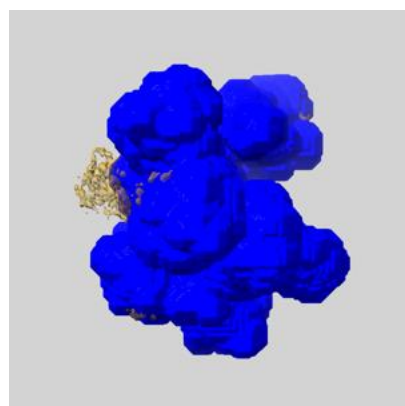
6.6 Mask visualisation [i](#)

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

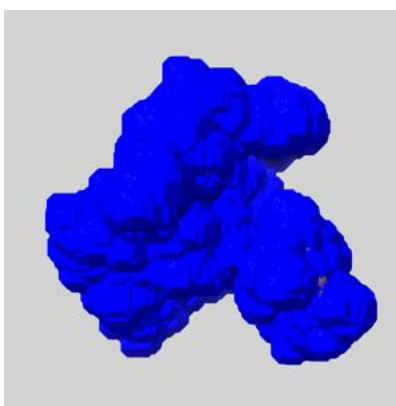
A mask typically either:

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

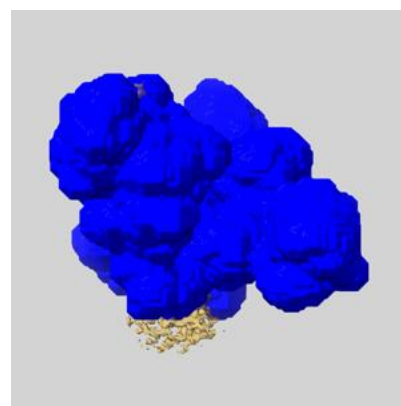
6.6.1 emd_55963_msk_1.map [i](#)



X



Y

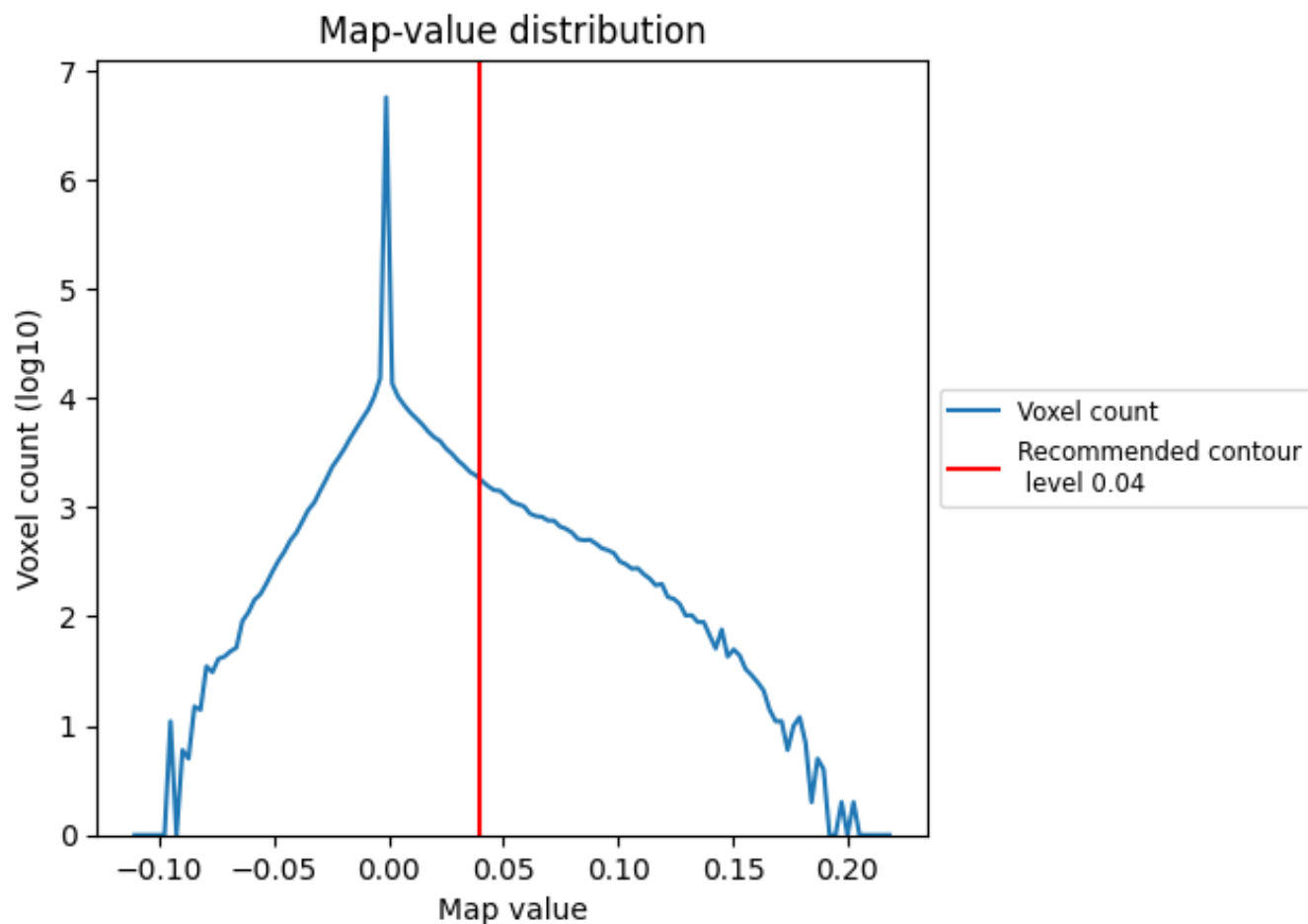


Z

7 Map analysis [i](#)

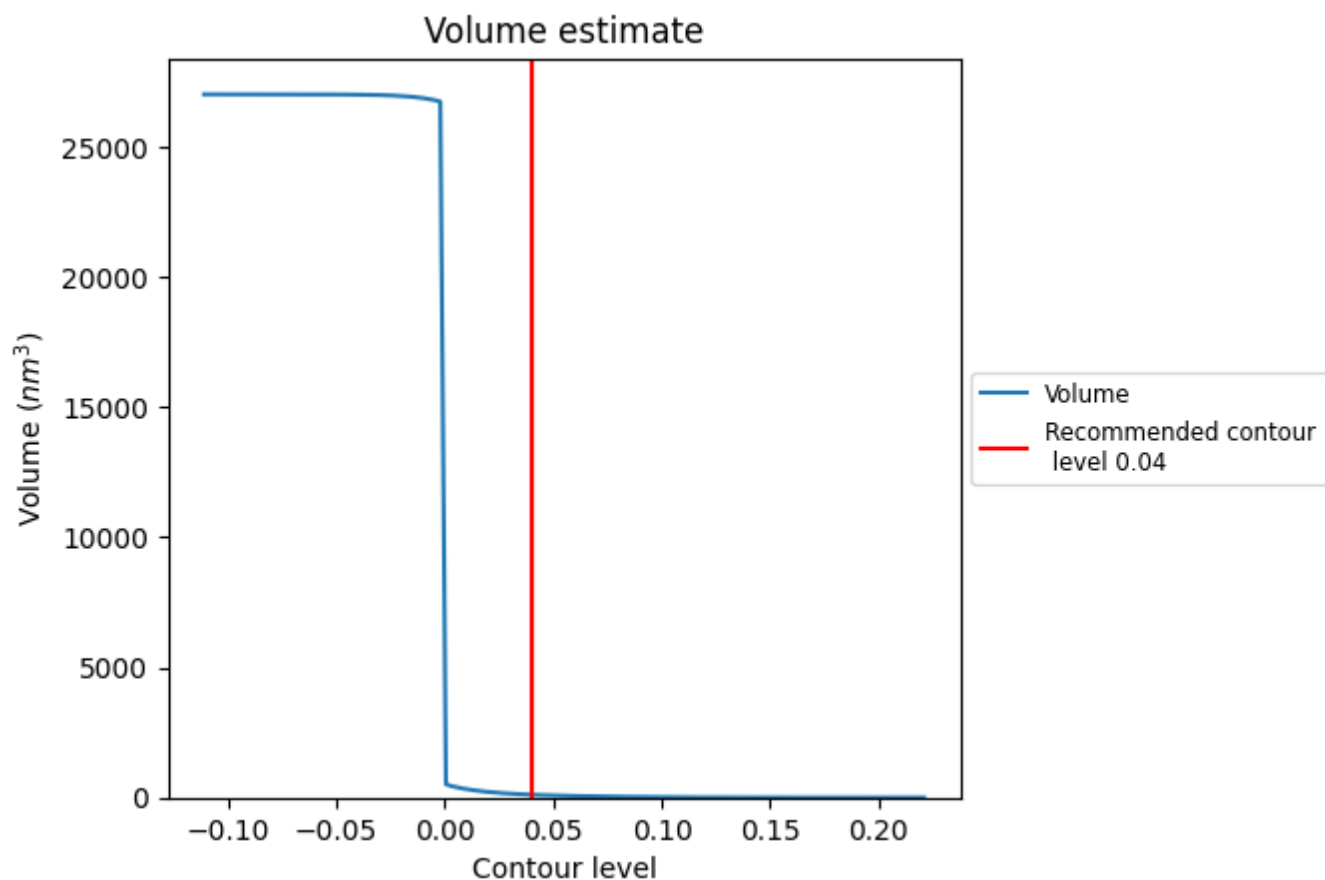
This section contains the results of statistical analysis of the map.

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.

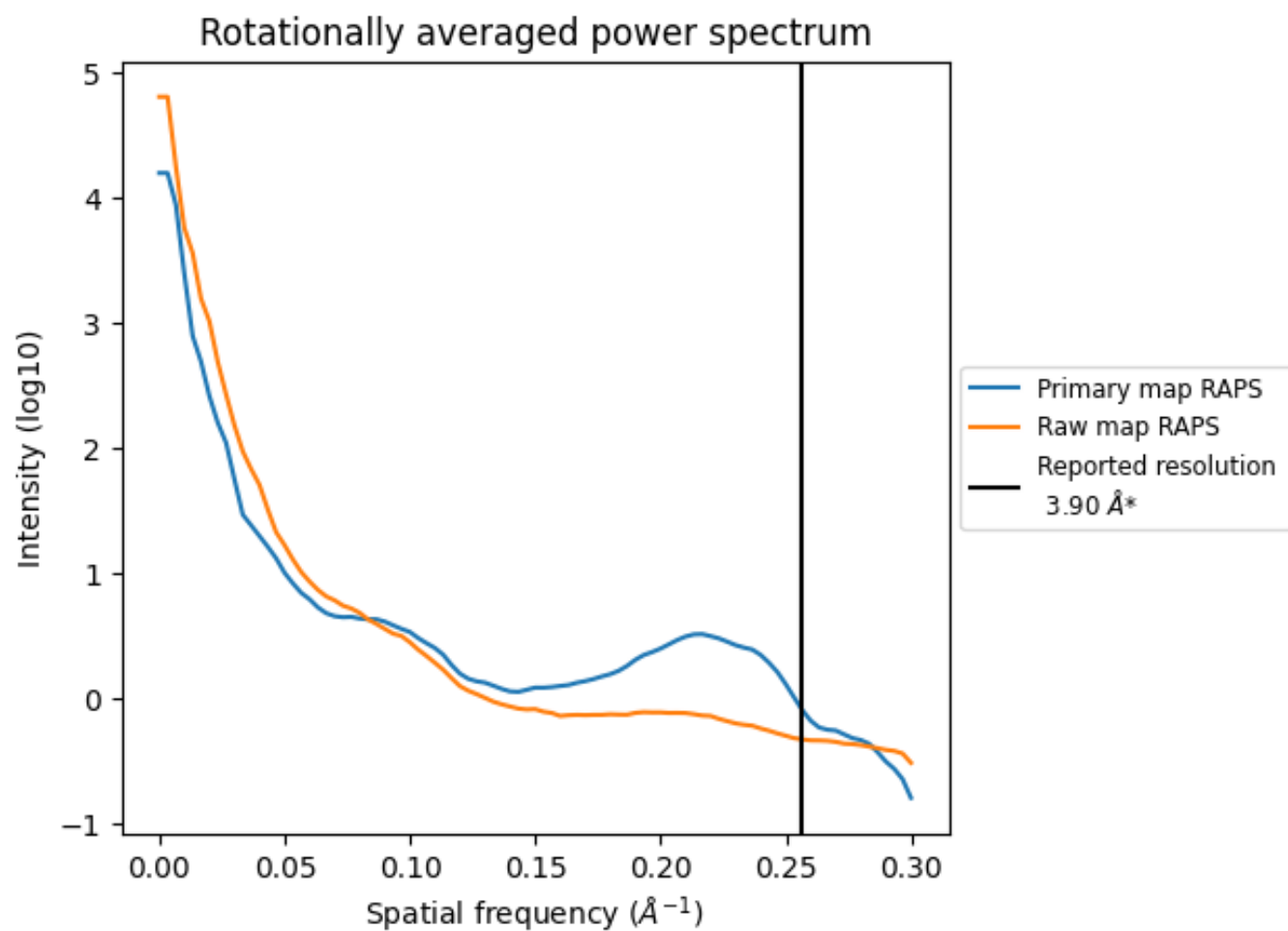
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 109 nm³; this corresponds to an approximate mass of 99 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 ⓘ

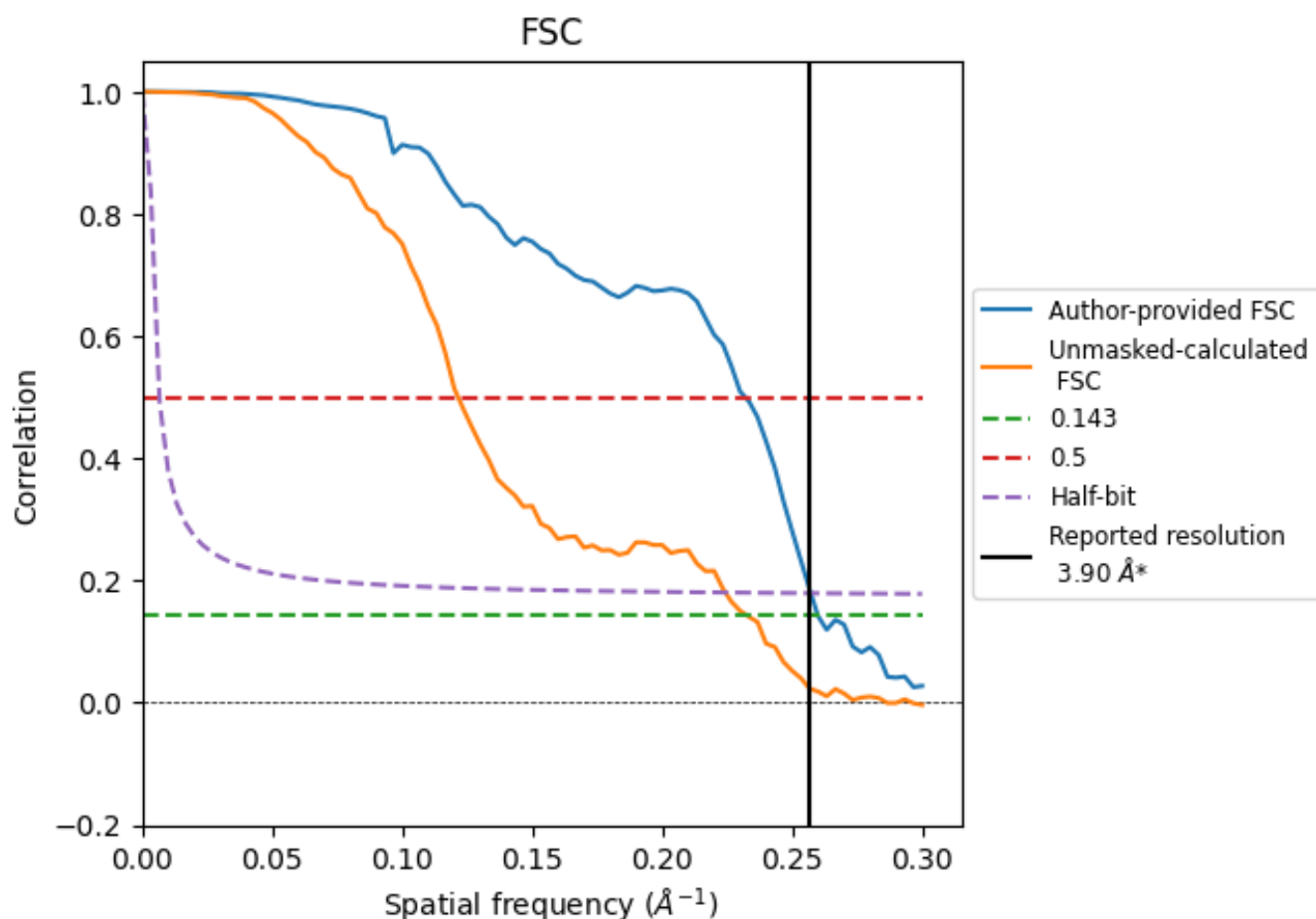


*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

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 0.256 Å⁻¹

8.2 Resolution estimates [i](#)

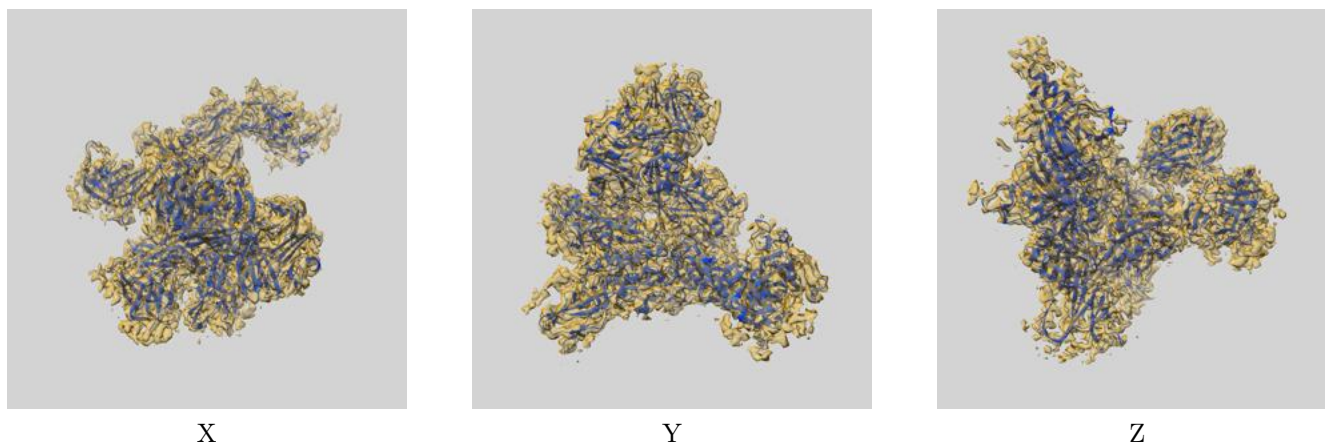
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	3.85	4.31	3.89
Unmasked-calculated*	4.30	8.24	4.46

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.30 differs from the reported value 3.9 by more than 10 %

9 Map-model fit [i](#)

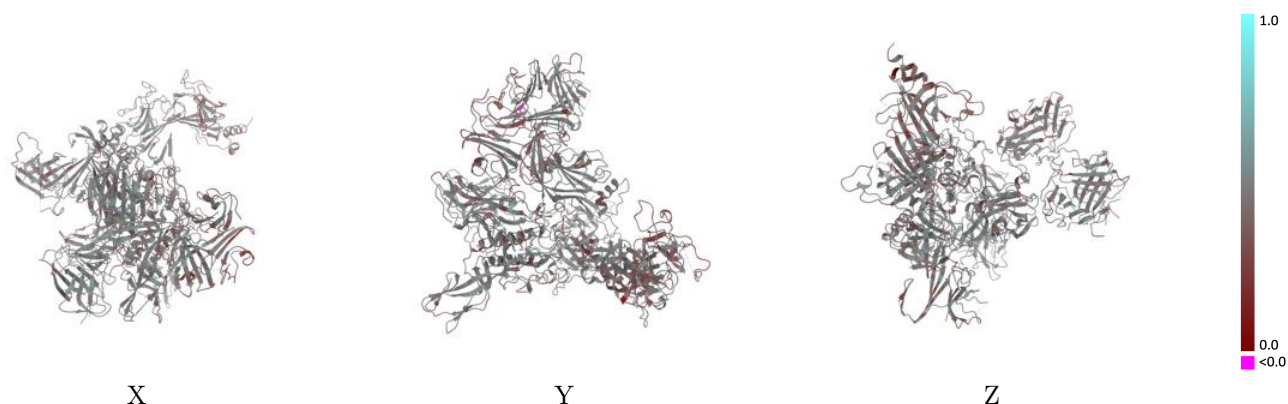
This section contains information regarding the fit between EMDB map EMD-55963 and PDB model 9TIO. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



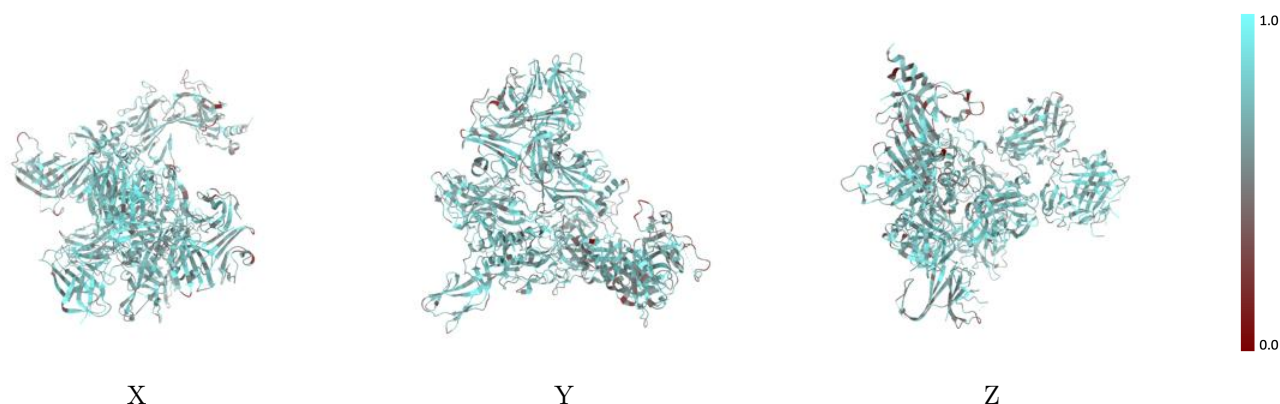
The images above show the 3D surface view of the map at the recommended contour level 0.04 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



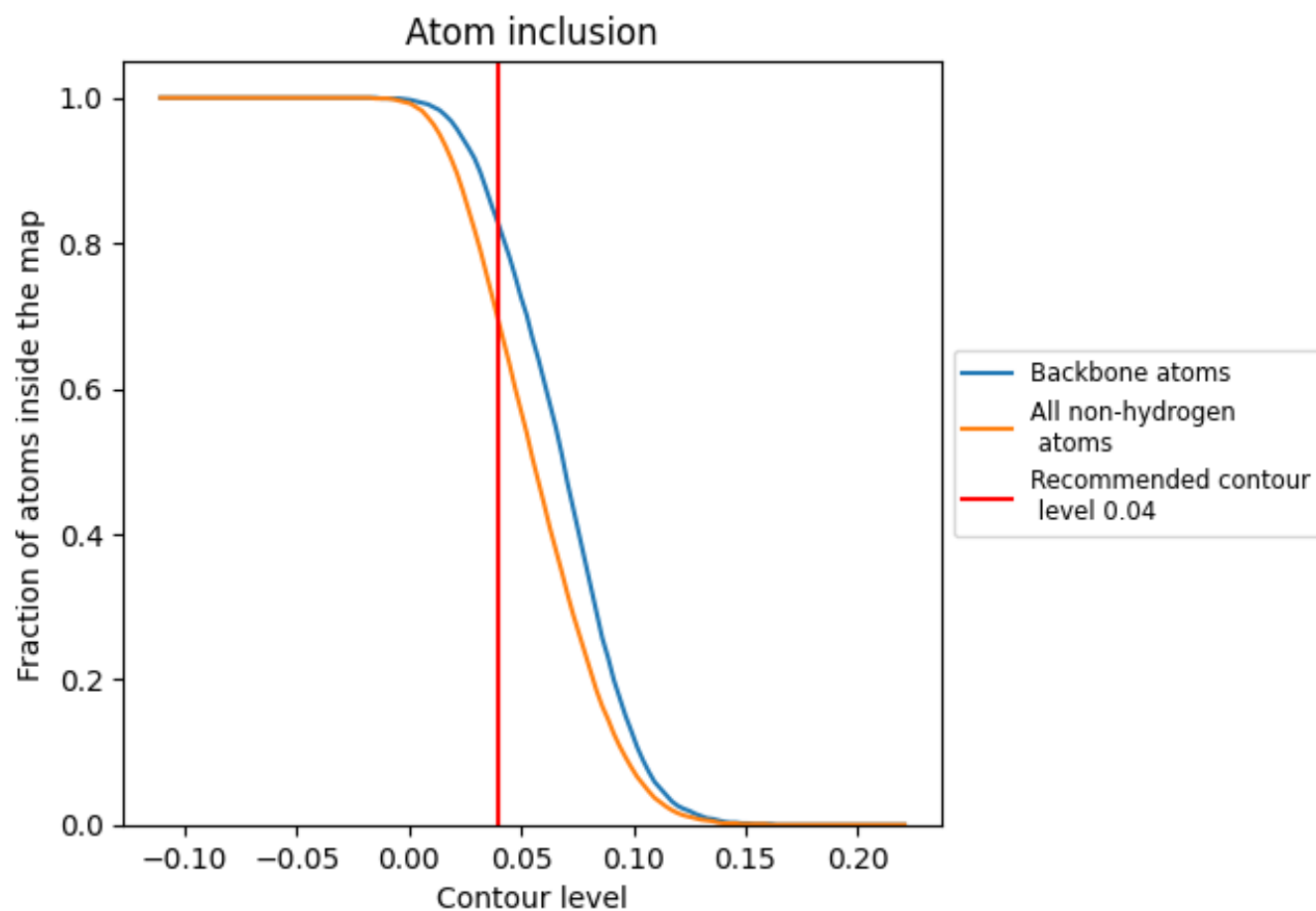
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.04).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 69% 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 (0.04) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6910	0.4440
A	0.7120	0.4670
B	0.6950	0.4550
C	0.6220	0.4270
D	0.7020	0.4600
E	0.6810	0.4400
F	0.7060	0.4500
G	0.7320	0.4620
H	0.6750	0.4120
I	0.7150	0.4500
J	0.6850	0.4250
K	0.5980	0.3920

1.0

0.0

<0.0