

wwPDB EM Validation Summary Report (i)

Feb 25, 2024 - 01:12 PM EST

PDB ID : 6XGQ EMDB ID EMD-22182 : Title YSD1 bacteriophage capsid : Authors Hardy, J.M.; Dunstan, R.; Venugopal, H.; Lithgow, T.J.; Coulibaly, F.J. : Deposited on 2020-06-17 : 3.80 Å(reported) Resolution : Based on initial model 6XGP •

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 (1)) were used in the production of this report:

EMDB validation analysis	:	0.0.1.dev70
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 3.80 Å.

Ramachandran outliers

Sidechain outliers

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	Percentile Ranks	Value
Ramachandran outliers		0
Sidechain outliers		0
Worse		Better
Percentile rel	ative to all structures	
Percentile rel	ative to all EM structures	
Metric	Whole archive (#Entries)	EM structures (#Entries)
	(#Entries)	(#Entries)

154571

154315

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 <=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.

4023

3826

Mol	Chain	Length	Quality of chain	
1	А	354	• 100%	
1	В	354	• 100%	
1	С	354	100%	
1	D	354	• 100%	
1	Е	354	100%	
1	F	354	• 100%	
1	G	354	94%	6%
2	a	139	89%	11%
2	b	139	89%	11%

Continued on next page...



		<i>i</i> previous		
Mol	Chain	Length	Quality of chain	
-		100		_
2	С	139	89%	11%
_	-			
2	d	139	89%	11%
2	е	139	89%	11%
2	f	139	89%	11%
2	g	139	89%	11%

Continued from previous page...



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 25706 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.

Mol	Chain	Residues		At	oms			AltConf	Trace
1	А	353	Total	С	Ν	0	S	0	0
	A	000	2791	1776	480	516	19	0	0
1	В	353	Total	С	Ν	0	S	0	0
	D	000	2791	1776	480	516	19	0	0
1	С	353	Total	С	Ν	0	S	0	0
	U	000	2791	1776	480	516	19	0	U
1	D	353	Total	С	Ν	0	S	0	0
	D	000	2791	1776	480	516	19	0	0
1	Е	353	Total	С	Ν	0	S	0	0
	Ľ	000	2791	1776	480	516	19	0	0
1	F	353	Total	С	Ν	0	S	0	0
	Г	000	2791	1776	480	516	19	0	0
1	G	332	Total	С	Ν	0	S	0	0
	G	552	2627	1675	451	483	18	0	0

• Molecule 1 is a protein called YSD1_17.

• Molecule 2 is a protein called YSD1_16.

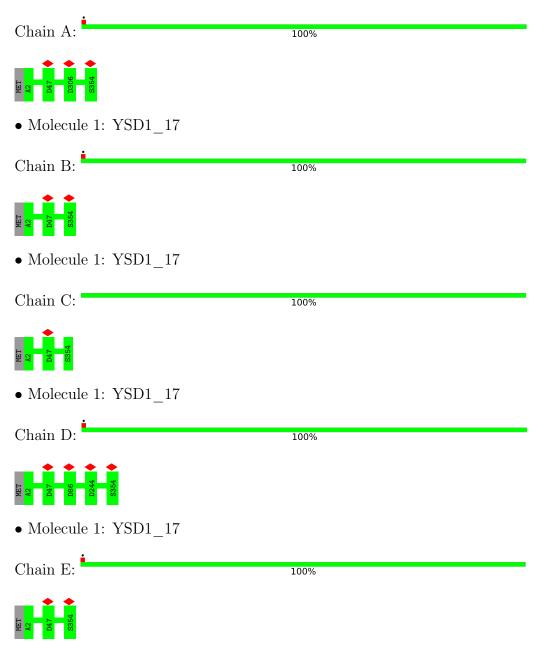
Mol	Chain	Residues		At	oms			AltConf	Trace
2	9	124	Total	С	Ν	0	S	0	0
	a	124	905	574	156	174	1	0	0
2	b	124	Total	С	Ν	Ο	\mathbf{S}	0	0
	U	124	905	574	156	174	1	0	0
2	0	124	Total	С	Ν	0	\mathbf{S}	0	0
	с	124	905	574	156	174	1	0	
2	d	124	Total	С	Ν	0	S	0	0
	u	124	905	574	156	174	1		0
2	0	124	Total	С	Ν	0	S	0	0
	е	124	905	574	156	174	1	0	0
2	f	124	Total	С	Ν	0	S	0	0
	1	124	905	574	156	174	1	0	0
2	ď	124	Total	С	Ν	0	S	0	0
	g	124	903	573	156	173	1	0	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.

 \bullet Molecule 1: YSD1_17





• Molecule 1: YSD1_17		
Chain F:	100%	
MET A2 B86 A309 A309 C310 C310 C310		
• Molecule 1: YSD1_17		
Chain G:	94%	6%
MET ALA GLY GLY L4 A170 A170 C217 L212 L120 L120 L120 L120 L120 L120 ASP ASP	clr hEU MET ASP SER THR ILE ARC GLU THR CLU THR CLU THR D243 A257 A257 S354	
• Molecule 2: YSD1_16		
Chain a:	89%	11%
MET ASN LLEU LLEU LLEU MET MET MET ALLA ALLA ALLA ALLA ALLA CLY TTO CLY THR SER SER SER SER AST		
• Molecule 2: YSD1_16		
Chain b:	89%	11%
MET ASN LEU LEU LEU MET MET ALA ALA ALA CLY THR THR THR FRO SER SER ASF		
• Molecule 2: YSD1_16		
Chain c:	89%	11%
MET LEU LEU LEU LEU LEU MET ALA ALA ALA TIA TIA TIA FRG SER FRG FRG FRG	<mark>, 139</mark>	
• Molecule 2: YSD1_16		
Chain d:	89%	11%
MET ASN LEU LEU LEU TER MET MET ALA ALA ALA ALA ALA ALA ALA ALA ALA ASP SFR SFR ASP		
• Molecule 2: YSD1_16		
Chain e:	89%	11%
MET ASN LEEU LEEU THR MET MET MET THR THR THR THR SER SER SER SER	<mark>. 139</mark>	





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	5449	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 $(e^-/\text{\AA}^2)$	27.24	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.080	Depositor
Minimum map value	-0.037	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	804.0, 804.0, 804.0	wwPDB
Map dimensions	600, 600, 600	wwPDB
Map angles $(^{\circ})$	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.34, 1.34, 1.34	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		
	Ullaill	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.46	0/2854	0.56	0/3866	
1	В	0.45	0/2854	0.57	0/3866	
1	С	0.45	0/2854	0.57	0/3866	
1	D	0.46	0/2854	0.56	0/3866	
1	Е	0.46	0/2854	0.57	0/3866	
1	F	0.46	0/2854	0.57	0/3866	
1	G	0.43	0/2687	0.58	0/3644	
2	a	0.43	0/927	0.57	0/1269	
2	b	0.45	0/927	0.58	0/1269	
2	с	0.45	0/927	0.56	0/1269	
2	d	0.46	0/927	0.58	0/1269	
2	е	0.45	0/927	0.58	0/1269	
2	f	0.46	0/927	0.59	0/1269	
2	g	0.44	0/925	0.57	0/1266	
All	All	0.45	0/26298	0.57	0/35720	

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)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percer	ntiles
1	А	351/354~(99%)	327~(93%)	24 (7%)	0	100	100
1	В	351/354~(99%)	333~(95%)	18 (5%)	0	100	100
1	С	351/354~(99%)	335~(95%)	16 (5%)	0	100	100
1	D	351/354~(99%)	333~(95%)	18 (5%)	0	100	100
1	Е	351/354~(99%)	325~(93%)	26 (7%)	0	100	100
1	F	351/354~(99%)	326 (93%)	25 (7%)	0	100	100
1	G	328/354~(93%)	305~(93%)	23 (7%)	0	100	100
2	a	120/139~(86%)	105 (88%)	15 (12%)	0	100	100
2	b	120/139~(86%)	109 (91%)	11 (9%)	0	100	100
2	с	120/139~(86%)	114 (95%)	6 (5%)	0	100	100
2	d	120/139~(86%)	111 (92%)	9 (8%)	0	100	100
2	е	120/139~(86%)	111 (92%)	9 (8%)	0	100	100
2	f	120/139~(86%)	108 (90%)	12 (10%)	0	100	100
2	g	120/139~(86%)	111 (92%)	9 (8%)	0	100	100
All	All	3274/3451~(95%)	3053~(93%)	221 (7%)	0	100	100

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

There are no Ramachandran outliers to report.

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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percer	ntiles
1	А	292/294~(99%)	292 (100%)	0	100	100
1	В	292/294~(99%)	292 (100%)	0	100	100
1	С	292/294~(99%)	292 (100%)	0	100	100
1	D	292/294~(99%)	292 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	Ε	292/294~(99%)	292 (100%)	0	100 100
1	\mathbf{F}	292/294~(99%)	292 (100%)	0	100 100
1	G	274/294~(93%)	274 (100%)	0	100 100
2	a	88/101 (87%)	88 (100%)	0	100 100
2	b	88/101~(87%)	88 (100%)	0	100 100
2	с	88/101 (87%)	88 (100%)	0	100 100
2	d	88/101 (87%)	88 (100%)	0	100 100
2	е	88/101 (87%)	88 (100%)	0	100 100
2	f	88/101 (87%)	88 (100%)	0	100 100
2	g	87/101~(86%)	87 (100%)	0	100 100
All	All	2641/2765~(96%)	2641 (100%)	0	100 100

Continued from previous page...

There are no protein residues with a non-rotameric sidechain to report.

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such side chains are listed below:

Mol	Chain	Res	Type
1	G	138	GLN

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 monosaccharides 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.



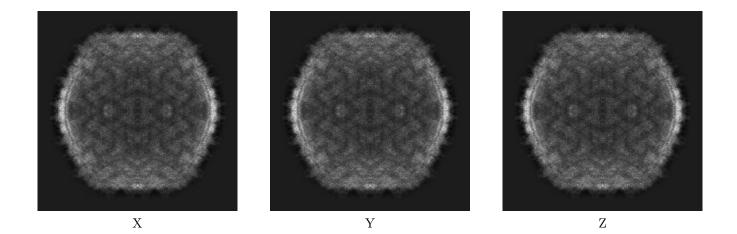
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-22182. These allow visual inspection of the internal detail of the map and identification of artifacts.

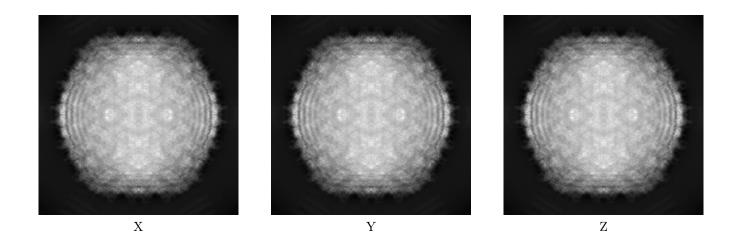
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

6.1.1 Primary map



6.1.2 Raw map



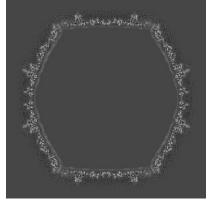
The images above show the map projected in three orthogonal directions.



6.2 Central slices (i)

6.2.1 Primary map





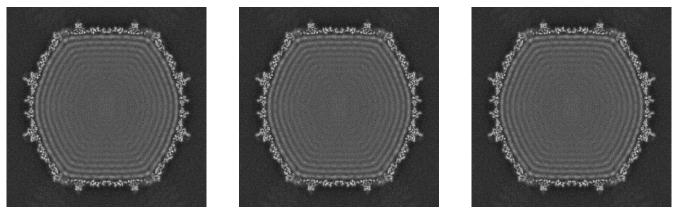


X Index: 300

Y Index: 300

Z Index: 300

6.2.2 Raw map



X Index: 300

Y Index: 300



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

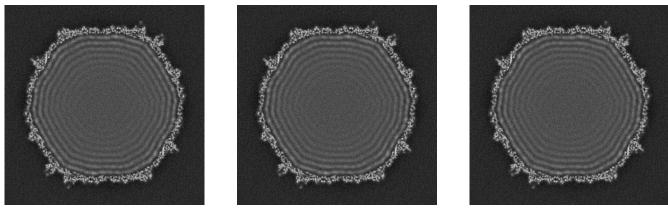


Y Index: 78



Z Index: 78

6.3.2 Raw map



X Index: 247

Y Index: 247

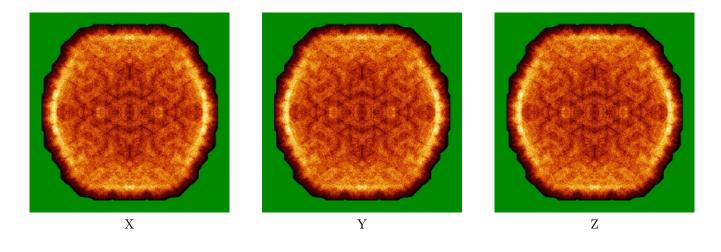


The images above show the largest variance slices of the map in three orthogonal directions.

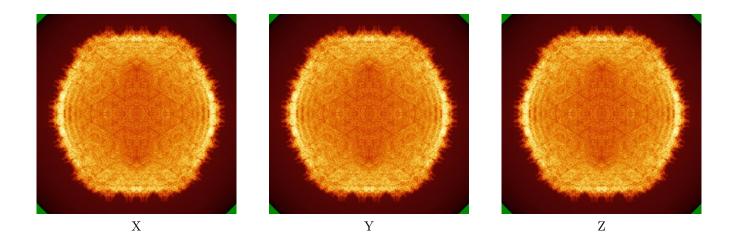


6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



6.4.2 Raw map

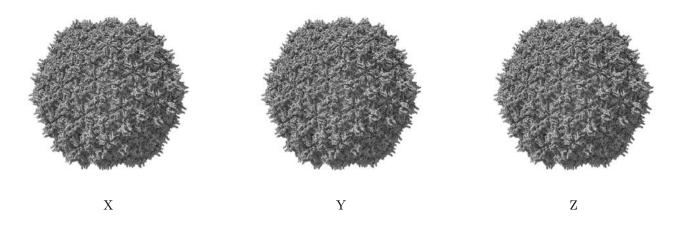


The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.



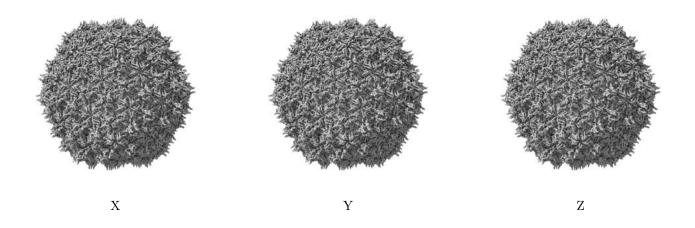
6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.015. 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.



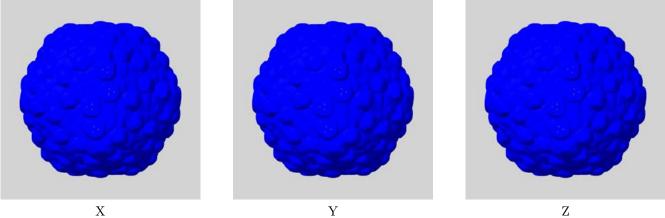
Mask visualisation (i) 6.6

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

$emd_{22182}msk_{1.map}$ (i) 6.6.1

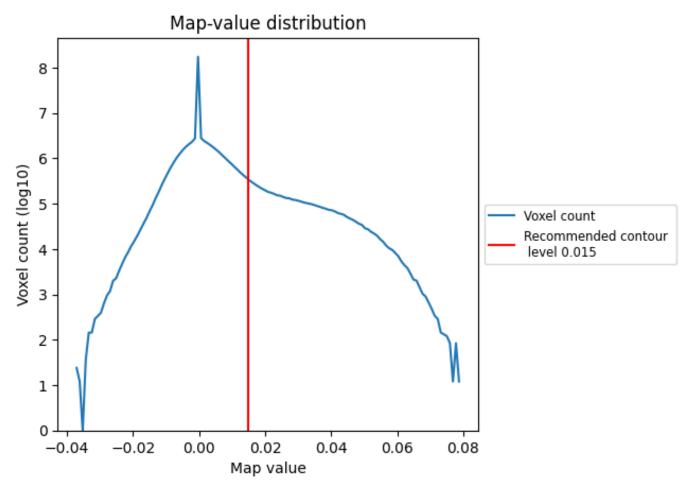




7 Map analysis (i)

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

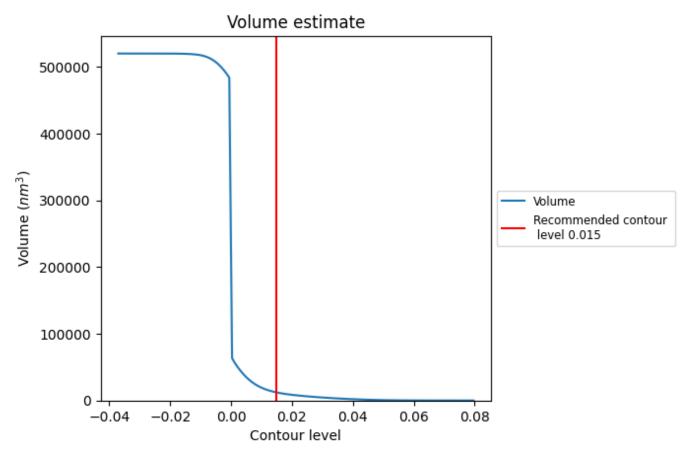
7.1 Map-value distribution (i)



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)

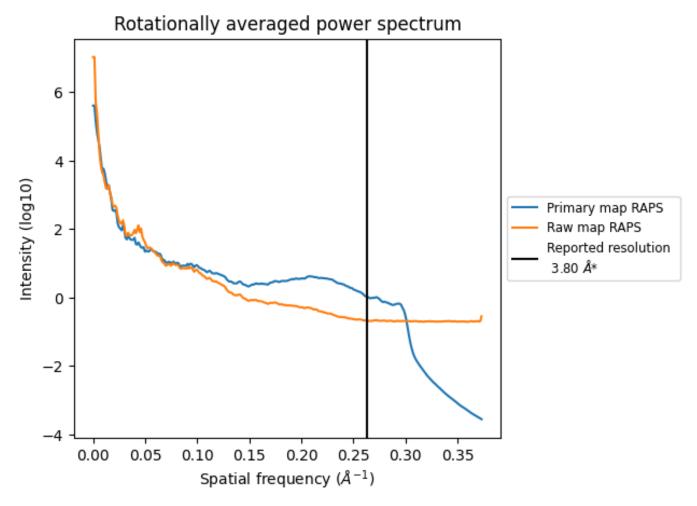


The volume at the recommended contour level is $12003~{\rm nm^3};$ this corresponds to an approximate mass of 10842 kDa.

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)



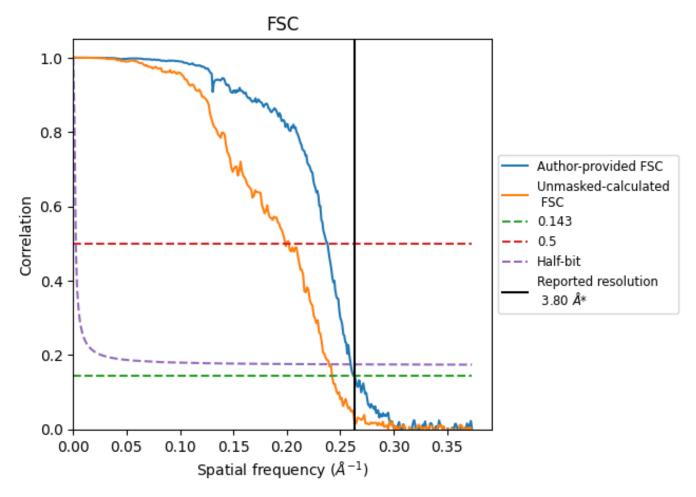
*Reported resolution corresponds to spatial frequency of 0.263 ${\rm \AA^{-1}}$



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 0.263 ${\rm \AA}^{-1}$



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	3.80	-	-
Author-provided FSC curve	3.80	4.20	3.85
Unmasked-calculated*	4.13	5.04	4.17

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

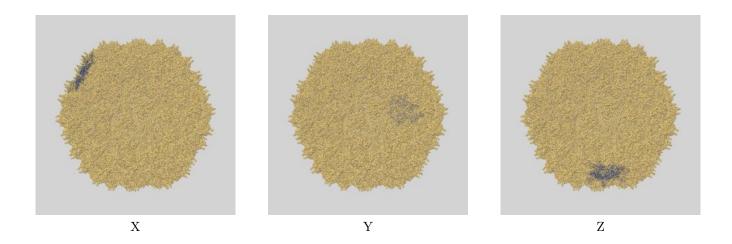


9 Map-model fit (i)

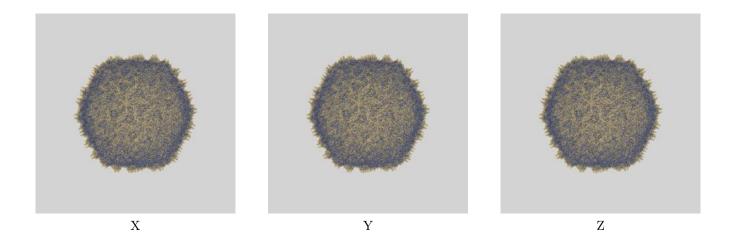
This section contains information regarding the fit between EMDB map EMD-22182 and PDB model 6XGQ. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlays

9.1.1 Map-model overlay (i)



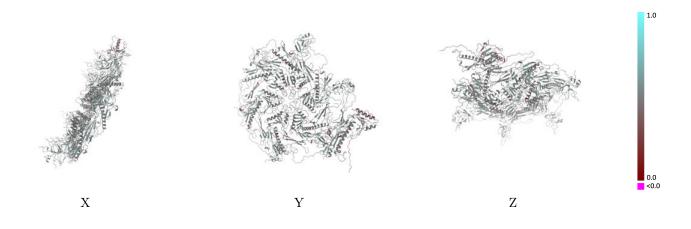
9.1.2 Map-model assembly overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.015 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.

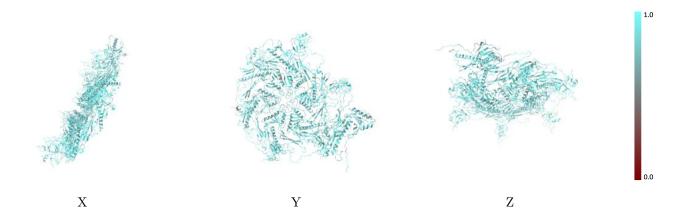


9.2 Q-score mapped to coordinate model (i)



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.

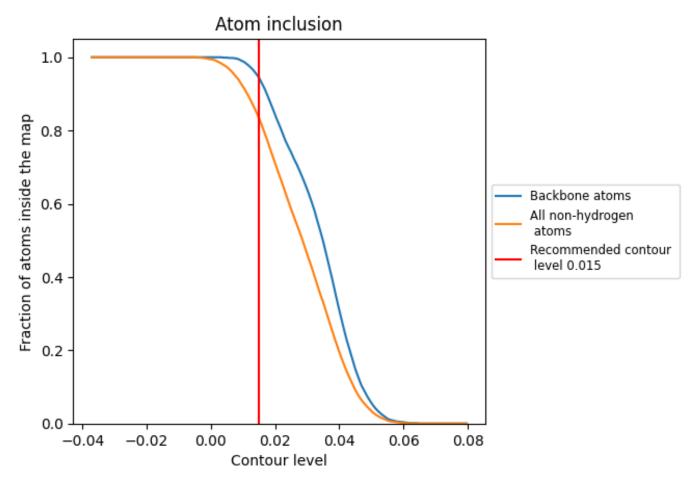
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.015).



9.4 Atom inclusion (i)



At the recommended contour level, 94% of all backbone atoms, 83% of all non-hydrogen atoms, are inside the map.



Map-model fit summary (i) 9.5

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	0.8330	0.4880	
А	0.8330	0.4920	1.0
В	0.8330	0.4890	
С	0.8340	0.4930	
D	0.8390	0.4920	
Е	0.8370	0.4930	
F	0.8280	0.4910	
G	0.7920	0.4610	
a	0.8350	0.4900	
b	0.8490	0.4840	
с	0.8430	0.4910	0.0
d	0.8510	0.4920	• <0.0
е	0.8510	0.4880	
f	0.8570	0.4930	
g	0.8520	0.4810	

