



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 03:45 AM UTC

PDB ID : 7KQE / pdb_00007kqe
EMDB ID : EMD-22997
Title : SARS-CoV-2 spike glycoprotein:Fab 3D11 complex
Authors : Asarnow, D.; Charles, C.; Cheng, Y.
Deposited on : 2020-11-15
Resolution : 2.88 Å(reported)
Based on initial model : 7A94

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.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

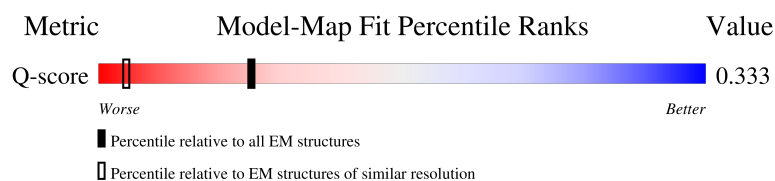
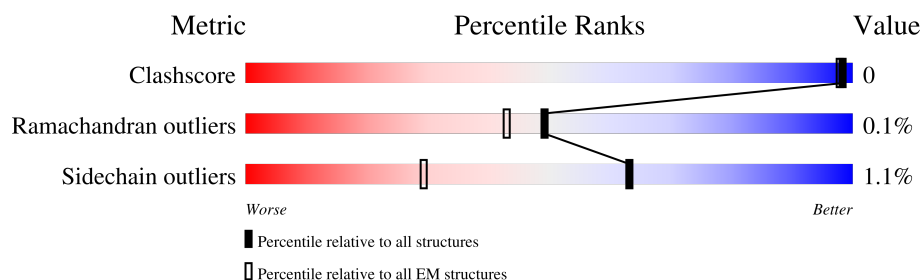
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.88 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.



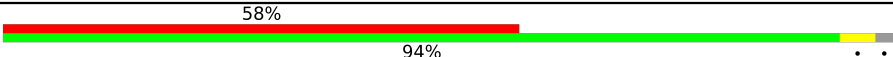
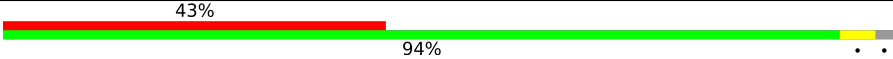
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	12111 (2.38 - 3.38)

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	H	448	47% 53%
1	I	448	19% 47% 53%
1	J	448	26% 47% 53%
2	L	216	11% 94% ..

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Mol	Chain	Length	Quality of chain
2	M	216	
2	N	216	
3	A	1208	
3	B	1208	
3	C	1208	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 68370 atoms, of which 33744 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 Fab 3D11 heavy chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	H	211	Total	C	H	N	O	S	0	0
			3187	1026	1580	271	306	4		
1	J	211	Total	C	H	N	O	S	0	0
			3187	1026	1580	271	306	4		
1	I	211	Total	C	H	N	O	S	0	0
			3187	1026	1580	271	306	4		

- Molecule 2 is a protein called Fab 3D11 light chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	L	212	Total	C	H	N	O	S	0	0
			3117	990	1526	266	330	5		
2	M	212	Total	C	H	N	O	S	0	0
			3117	990	1526	266	330	5		
2	N	212	Total	C	H	N	O	S	0	0
			3117	990	1526	266	330	5		

- Molecule 3 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	A	1066	Total	C	H	N	O	S	0	0
			16486	5324	8142	1394	1588	38		
3	B	1066	Total	C	H	N	O	S	0	0
			16486	5324	8142	1394	1588	38		
3	C	1066	Total	C	H	N	O	S	0	0
			16486	5324	8142	1394	1588	38		

There are 15 discrepancies between the modelled and reference sequences:

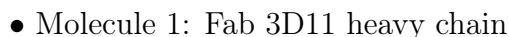
Chain	Residue	Modelled	Actual	Comment	Reference
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2
A	685	SER	ARG	conflict	UNP P0DTC2

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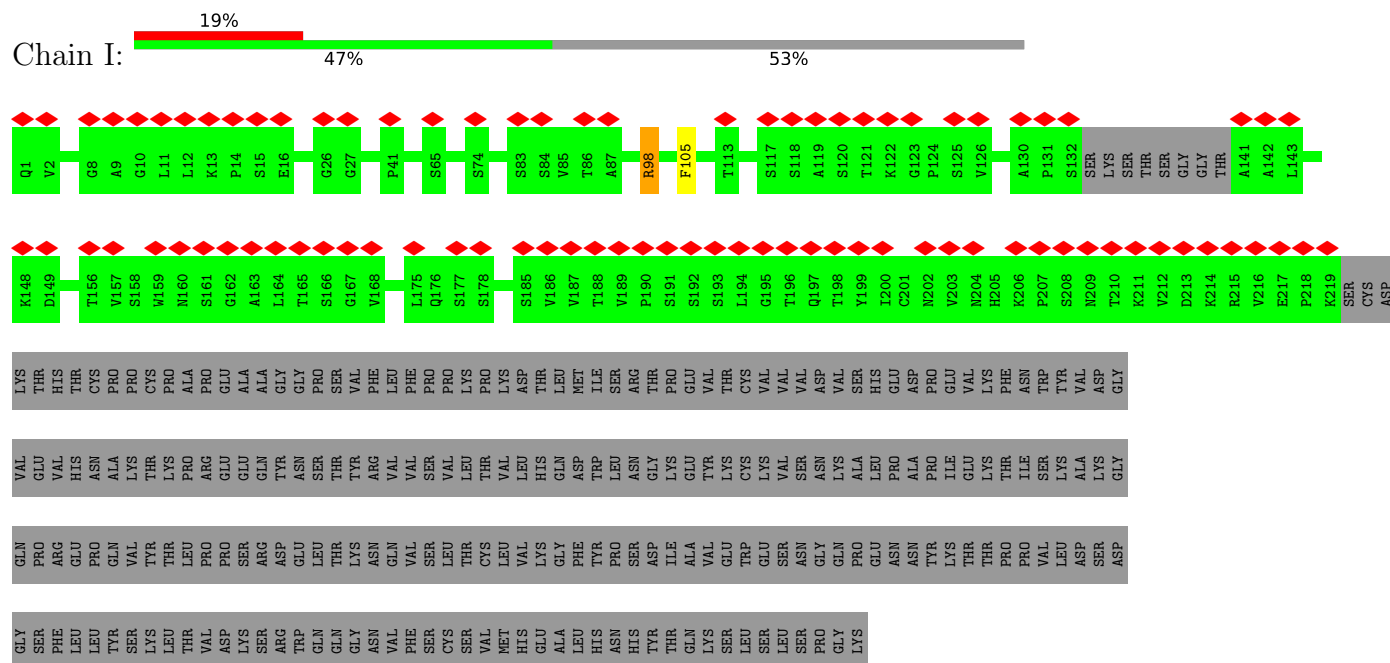
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Chain	Residue	Modelled	Actual	Comment	Reference
A	986	PRO	LYS	conflict	UNP P0DTC2
A	987	PRO	VAL	conflict	UNP P0DTC2
B	682	GLY	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
C	682	GLY	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	986	PRO	LYS	conflict	UNP P0DTC2
C	987	PRO	VAL	conflict	UNP P0DTC2

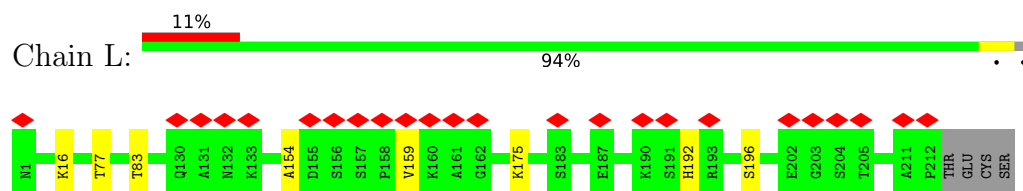
- Molecule 1: Fab 3D11 heavy chain



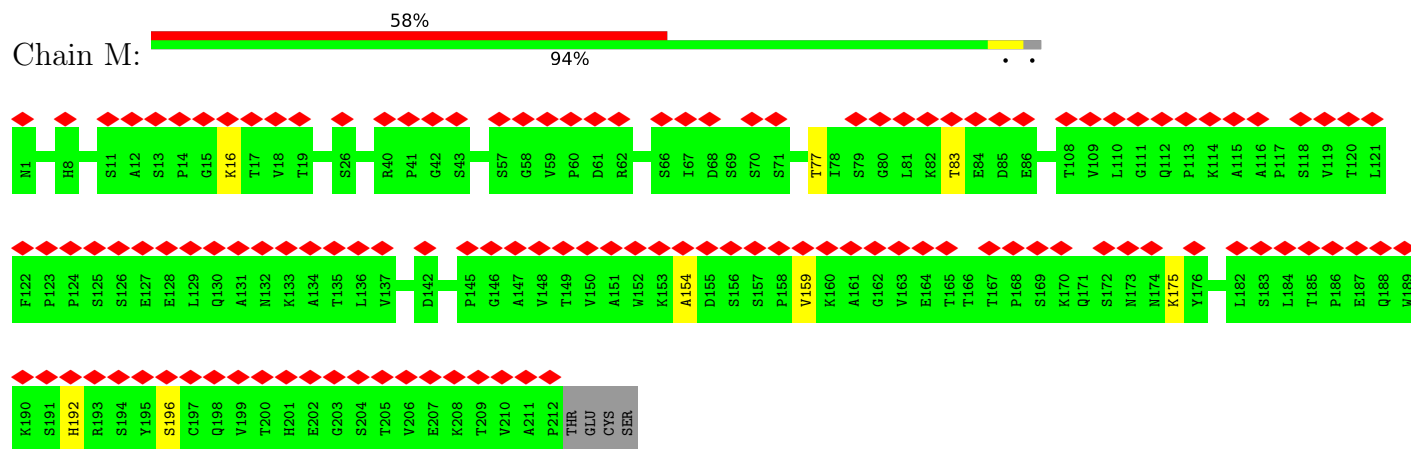
- Molecule 1: Fab 3D11 heavy chain



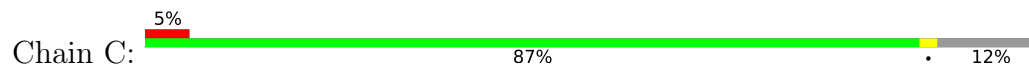
- Molecule 2: Fab 3D11 light chain

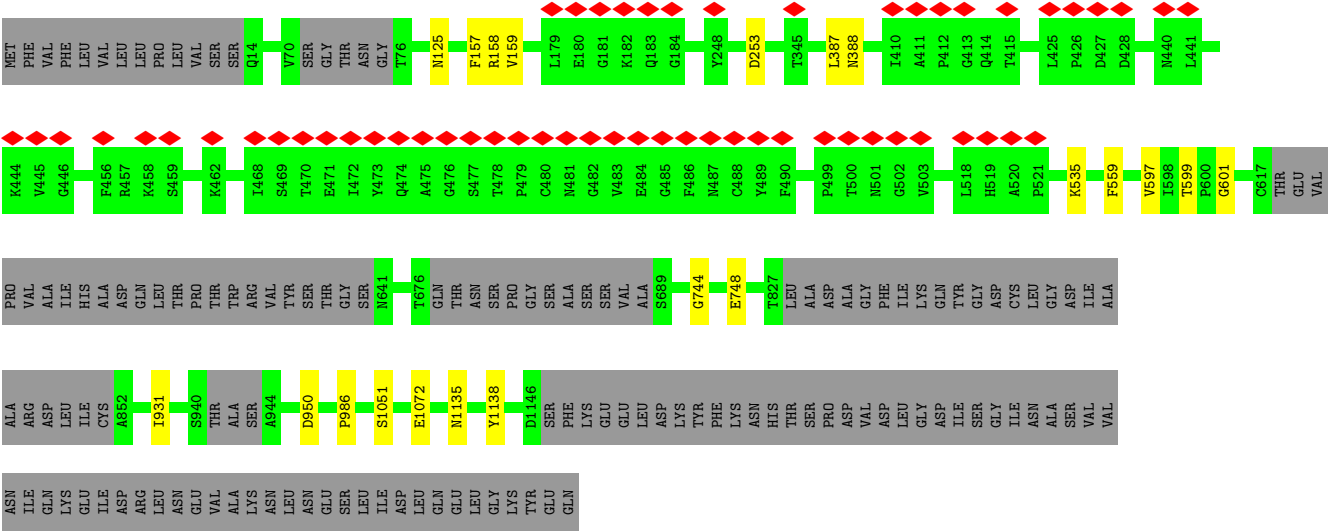


- Molecule 2: Fab 3D11 light chain



- Molecule 2: Fab 3D11 light chain





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	197150	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 ($e^-/\text{\AA}^2$)	67.7	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.291	Depositor
Minimum map value	-0.777	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.086	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	427.52, 427.52, 427.52	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4250666, 1.4250666, 1.4250666	Depositor

5 Model quality

5.1 Standard geometry

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	H	0.75	0/1652	1.30	0/2258
1	I	0.75	0/1652	1.30	0/2258
1	J	0.74	0/1652	1.29	0/2258
2	L	0.72	0/1630	1.25	0/2228
2	M	0.72	0/1630	1.25	0/2228
2	N	0.72	0/1630	1.25	0/2228
3	A	0.73	0/8538	1.23	10/11618 (0.1%)
3	B	0.73	0/8538	1.22	4/11618 (0.0%)
3	C	0.72	0/8538	1.23	7/11618 (0.1%)
All	All	0.73	0/35460	1.24	21/48312 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	H	0	1
1	I	0	1
1	J	0	1
3	A	0	3
3	B	0	2
All	All	0	8

There are no bond length outliers.

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	125	ASN	CA-CB-CG	8.07	120.67	112.60
3	A	80	ASP	CA-CB-CG	6.22	118.82	112.60
3	A	1083	HIS	O-C-N	-5.86	114.80	122.59
3	B	986	PRO	N-CA-C	5.79	117.77	110.70
3	A	986	PRO	N-CA-C	5.73	117.69	110.70

There are no chirality outliers.

5 of 8 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	246	ARG	Sidechain
3	A	655	HIS	Sidechain
3	A	79	PHE	Sidechain
3	B	246	ARG	Sidechain
1	H	98	ARG	Sidechain

5.2 Too-close contacts

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	H	1607	1580	1579	1	0
1	I	1607	1580	1579	1	0
1	J	1607	1580	1579	1	0
2	L	1591	1526	1526	2	0
2	M	1591	1526	1526	2	0
2	N	1591	1526	1526	2	0
3	A	8344	8142	8136	1	0
3	B	8344	8142	8136	4	0
3	C	8344	8142	8136	3	0
All	All	34626	33744	33723	17	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 0.

The worst 5 of 17 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:154:ALA:HB3	2:M:159:VAL:HG21	1.89	0.54
2:N:154:ALA:HB3	2:N:159:VAL:HG21	1.89	0.53
2:L:154:ALA:HB3	2:L:159:VAL:HG21	1.89	0.53
3:C:599:THR:HG22	3:C:601:GLY:H	1.77	0.50
1:J:98:ARG:HD3	1:J:105:PHE:CE1	2.47	0.50

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	H	207/448 (46%)	194 (94%)	13 (6%)	0	100	100
1	I	207/448 (46%)	194 (94%)	13 (6%)	0	100	100
1	J	207/448 (46%)	194 (94%)	13 (6%)	0	100	100
2	L	210/216 (97%)	201 (96%)	9 (4%)	0	100	100
2	M	210/216 (97%)	201 (96%)	9 (4%)	0	100	100
2	N	210/216 (97%)	201 (96%)	9 (4%)	0	100	100
3	A	1054/1208 (87%)	978 (93%)	74 (7%)	2 (0%)	43	70
3	B	1054/1208 (87%)	986 (94%)	67 (6%)	1 (0%)	48	74
3	C	1054/1208 (87%)	981 (93%)	71 (7%)	2 (0%)	43	70
All	All	4413/5616 (79%)	4130 (94%)	278 (6%)	5 (0%)	49	74

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	158	ARG
3	A	158	ARG
3	B	744	GLY
3	A	594	GLY
3	C	744	GLY

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	H	180/396 (46%)	180 (100%)	0	100	100
1	I	180/396 (46%)	180 (100%)	0	100	100
1	J	180/396 (46%)	180 (100%)	0	100	100
2	L	182/186 (98%)	177 (97%)	5 (3%)	39	70
2	M	182/186 (98%)	177 (97%)	5 (3%)	39	70
2	N	182/186 (98%)	177 (97%)	5 (3%)	39	70
3	A	932/1053 (88%)	921 (99%)	11 (1%)	63	84
3	B	932/1053 (88%)	924 (99%)	8 (1%)	70	88
3	C	932/1053 (88%)	924 (99%)	8 (1%)	70	88
All	All	3882/4905 (79%)	3840 (99%)	42 (1%)	63	86

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	597	VAL
2	M	175	LYS
3	C	1051	SER
2	M	16	LYS
2	N	16	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 84 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	239	GLN
3	C	1011	GLN
3	C	450	ASN
3	C	755	GLN
1	J	39	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [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.

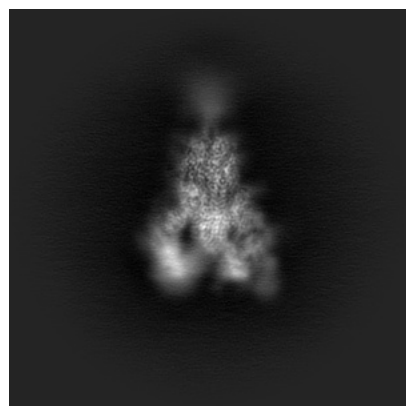
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22997. These allow visual inspection of the internal detail of the map and identification of artifacts.

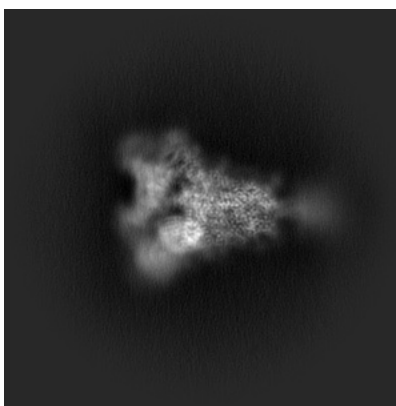
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

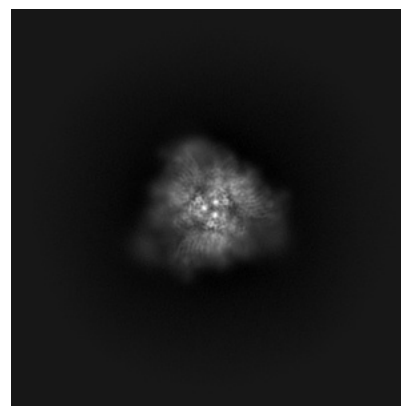
6.1.1 Primary map



X

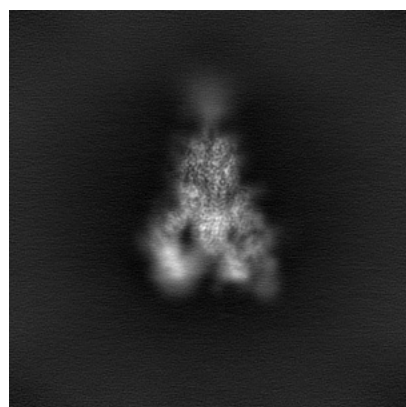


Y

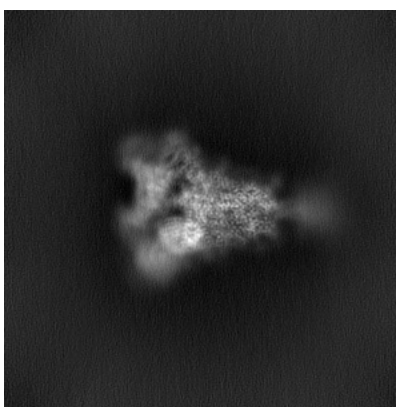


Z

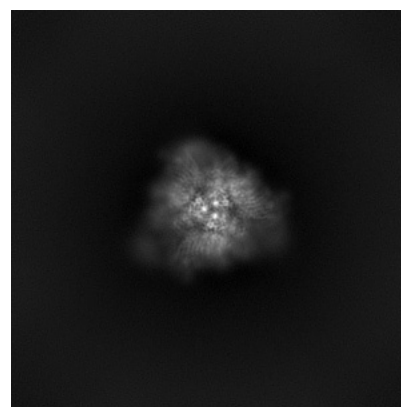
6.1.2 Raw map



X



Y

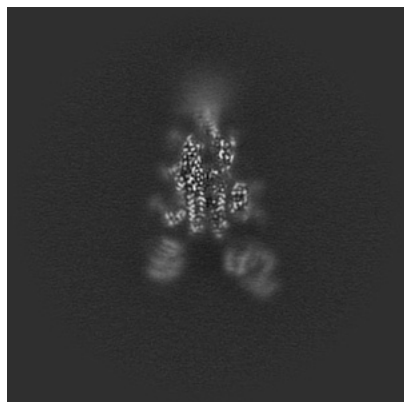


Z

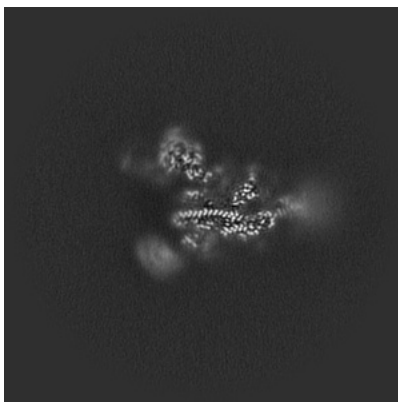
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

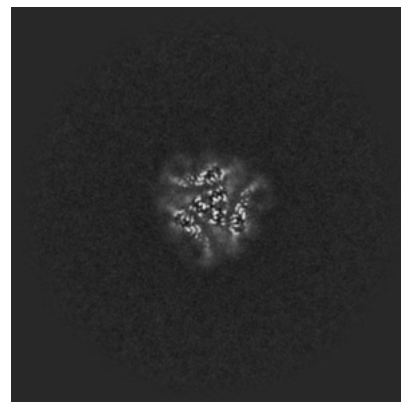
6.2.1 Primary map



X Index: 150

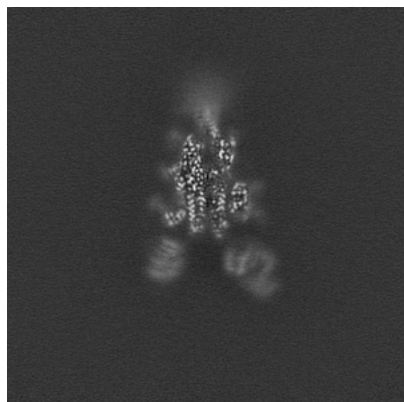


Y Index: 150



Z Index: 150

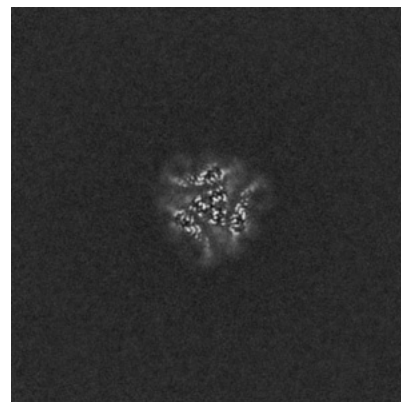
6.2.2 Raw map



X Index: 150



Y Index: 150



Z Index: 150

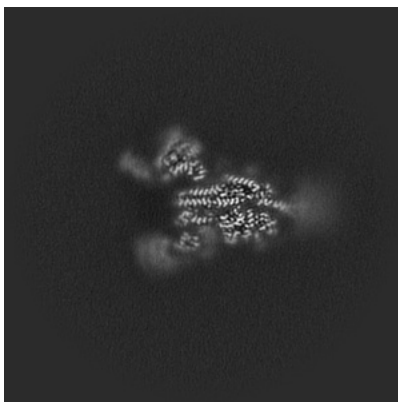
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

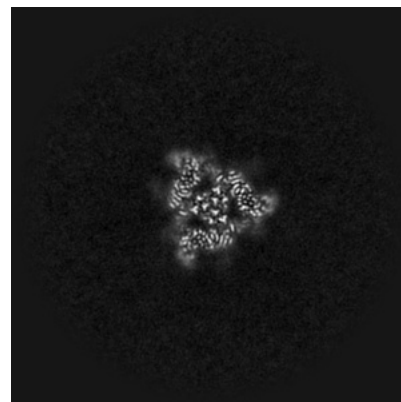
6.3.1 Primary map



X Index: 153



Y Index: 155

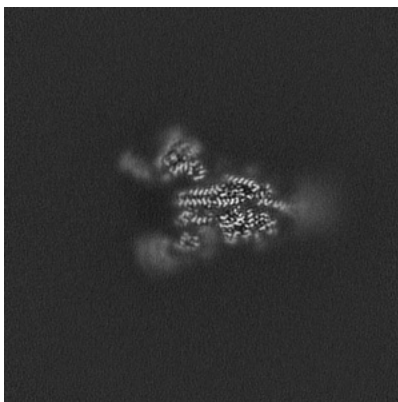


Z Index: 143

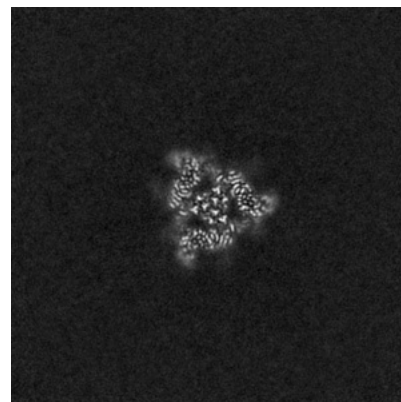
6.3.2 Raw map



X Index: 153



Y Index: 155

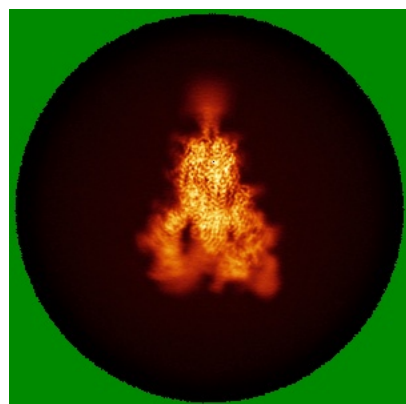


Z Index: 143

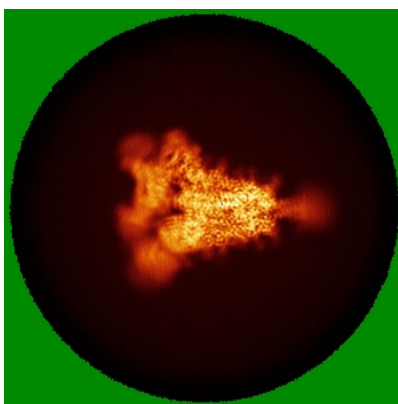
The images above show the largest variance slices of the map in three orthogonal directions.

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

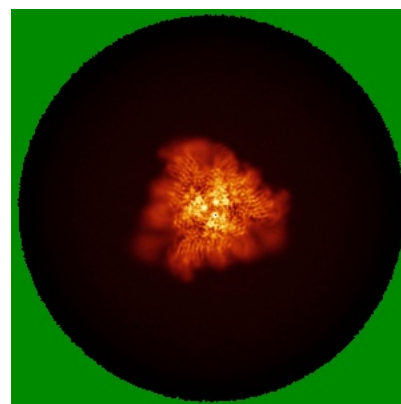
6.4.1 Primary map



X

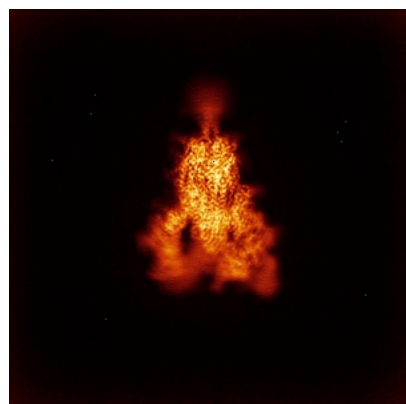


Y

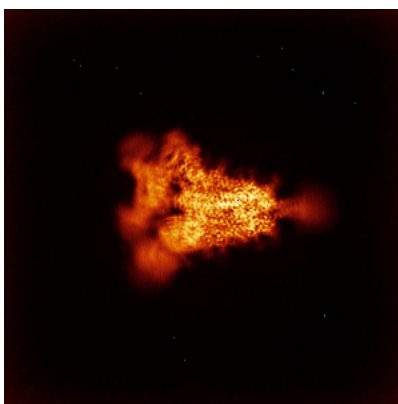


Z

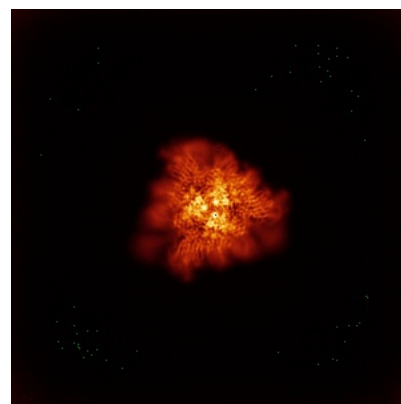
6.4.2 Raw map



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 0.5. 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.

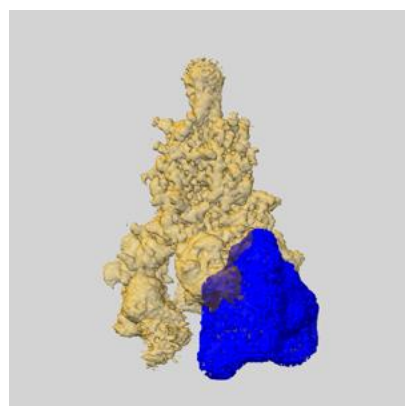
6.6 Mask visualisation [i](#)

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

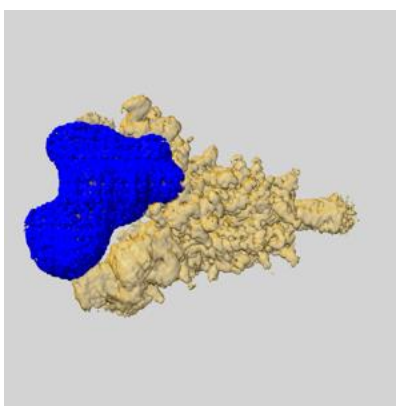
A mask typically either:

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

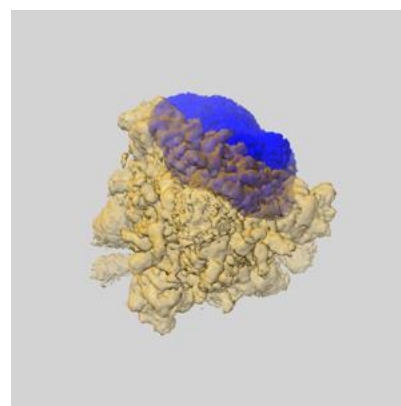
6.6.1 emd_22997_msk_1.map [i](#)



X



Y

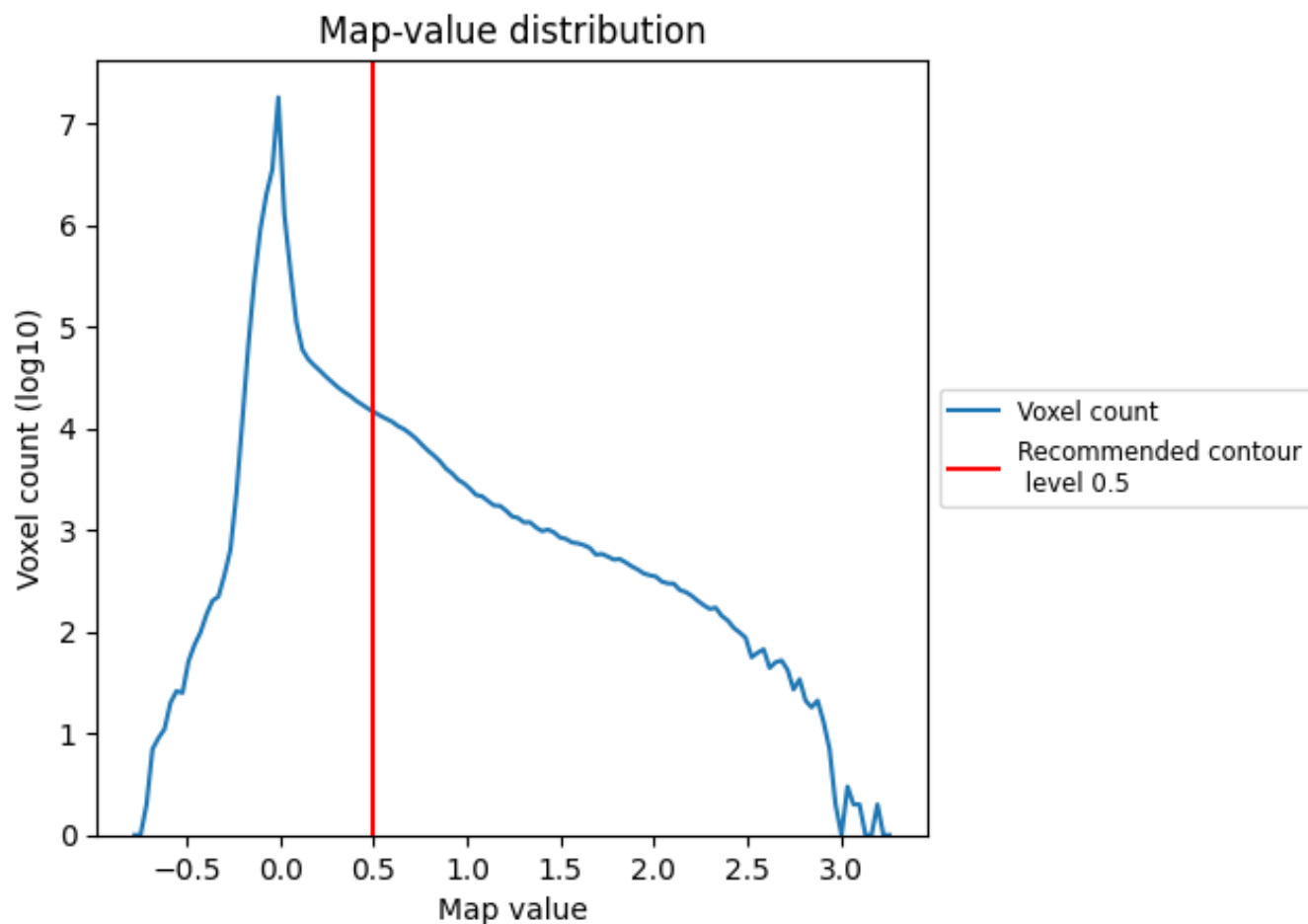


Z

7 Map analysis [i](#)

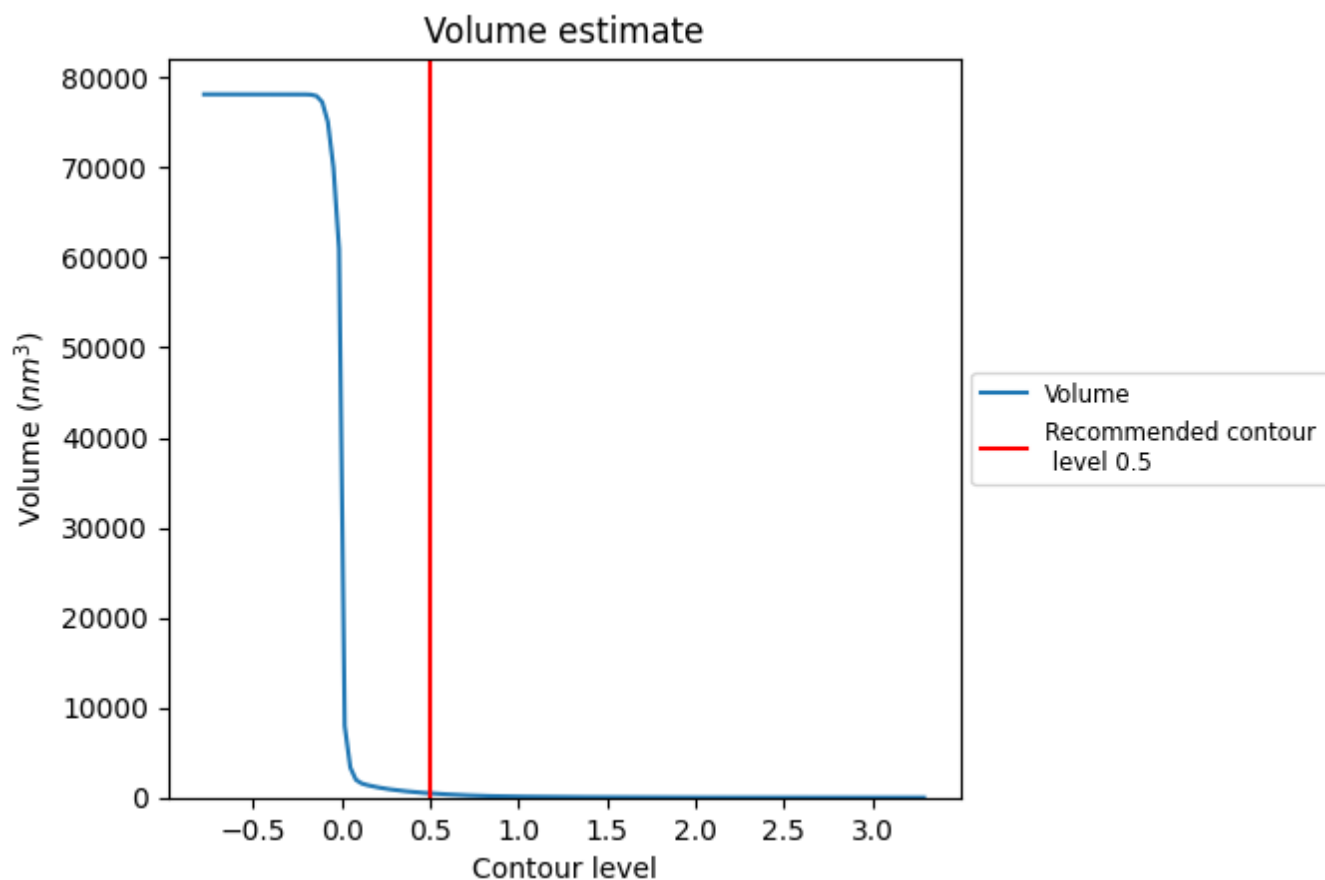
This section contains the results of statistical analysis of the map.

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.

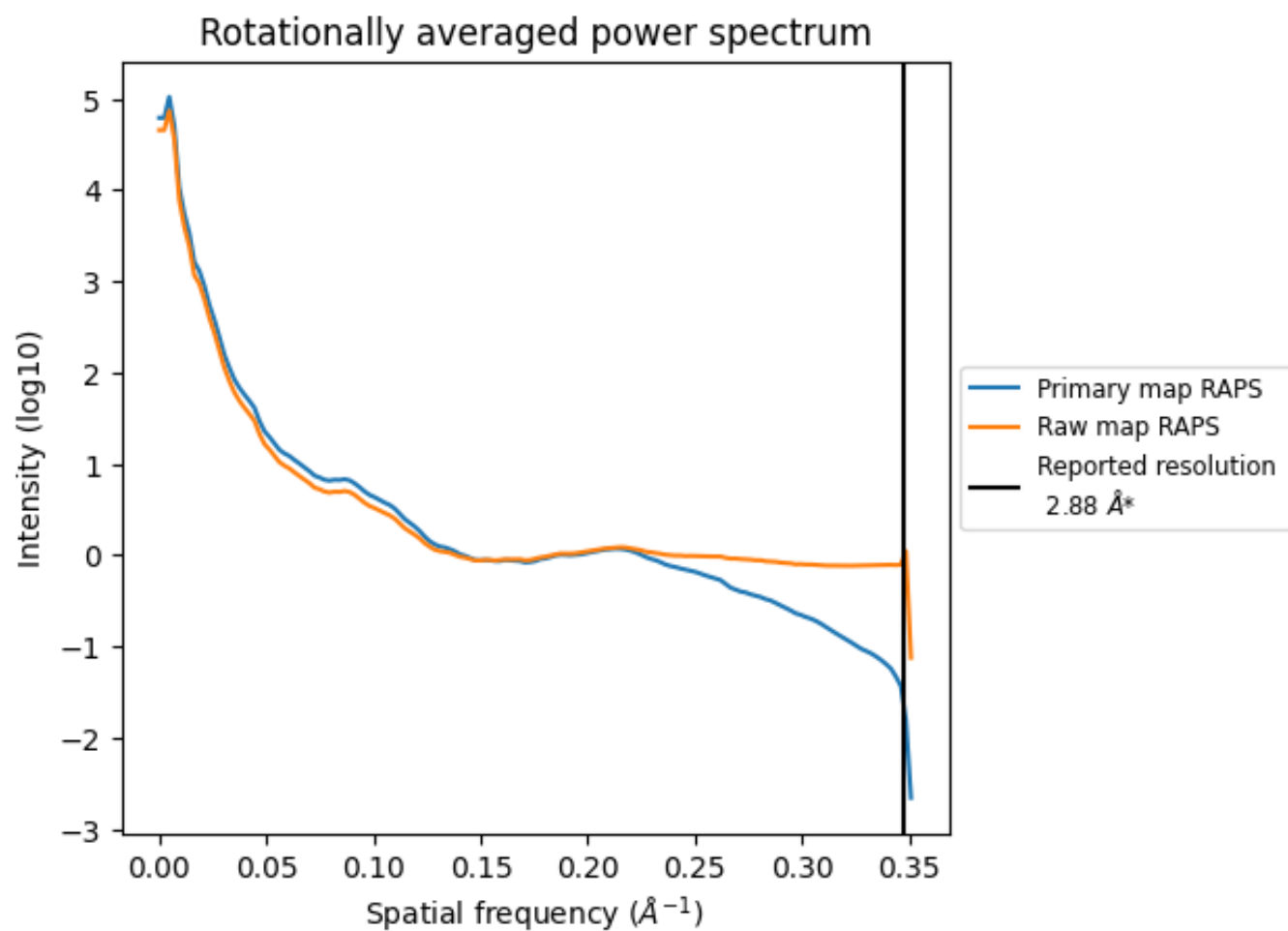
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 473 nm³; this corresponds to an approximate mass of 427 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 ⓘ

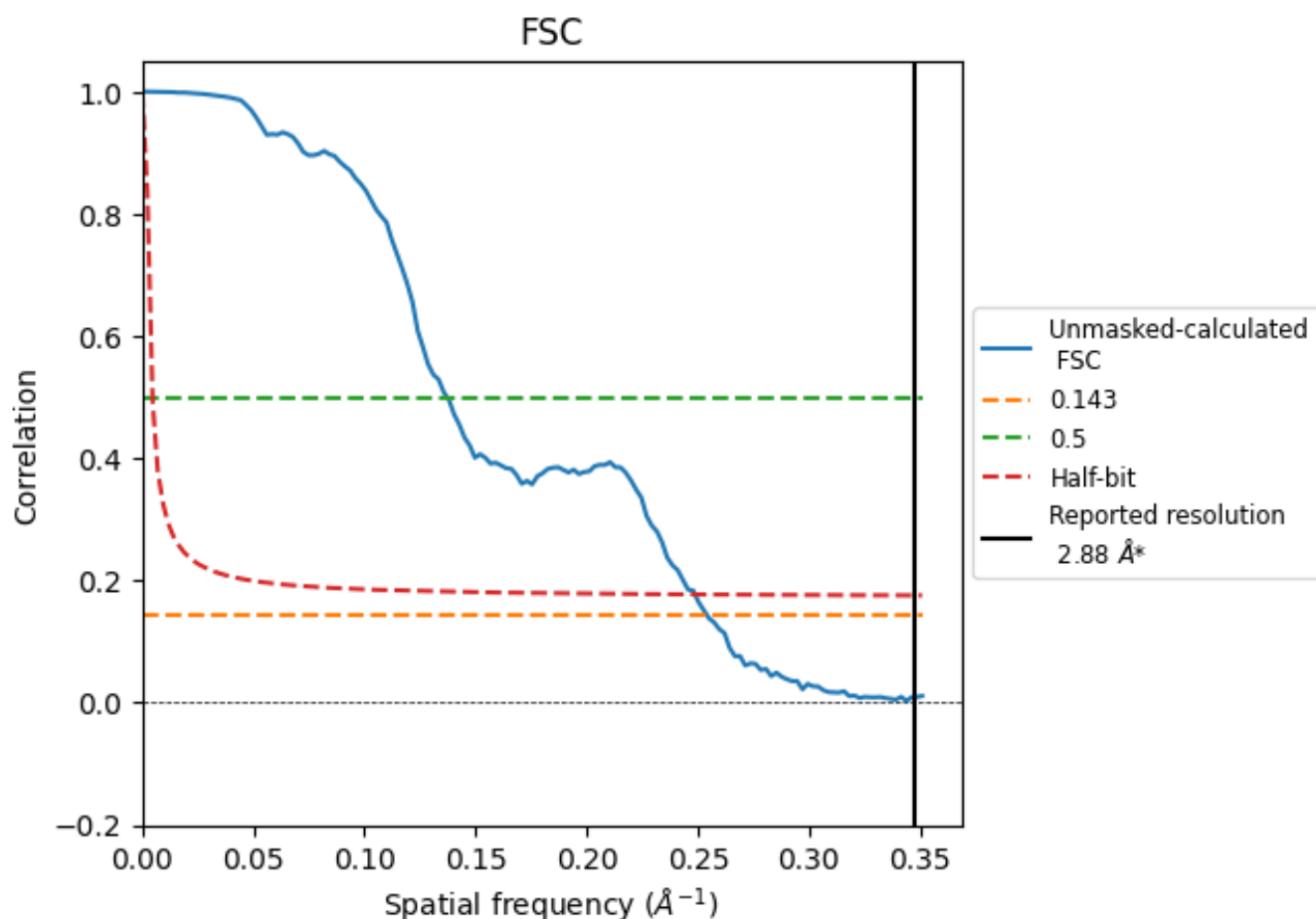


*Reported resolution corresponds to spatial frequency of 0.347 Å⁻¹

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 0.347 Å⁻¹

8.2 Resolution estimates [i](#)

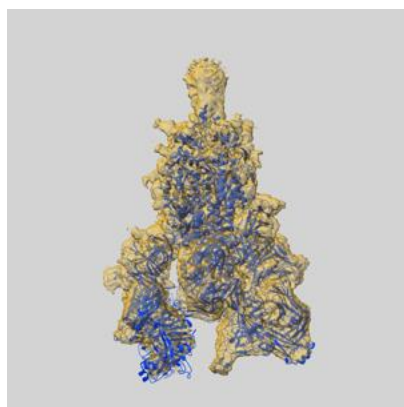
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.88	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.94	7.29	4.02

*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 3.94 differs from the reported value 2.88 by more than 10 %

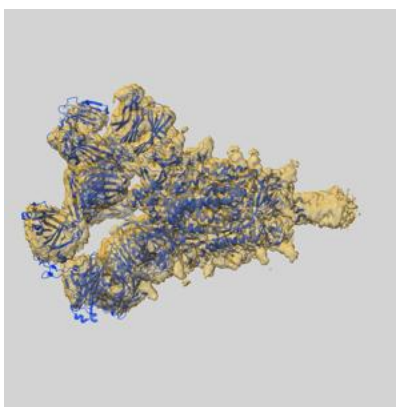
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22997 and PDB model 7KQE. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

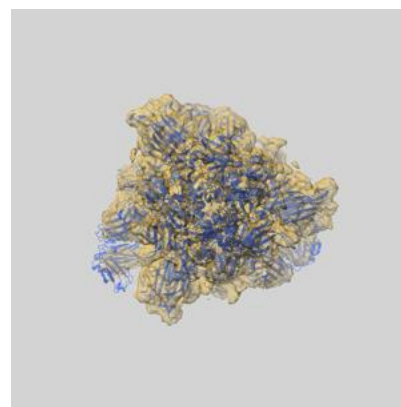
9.1 Map-model overlay [i](#)



X



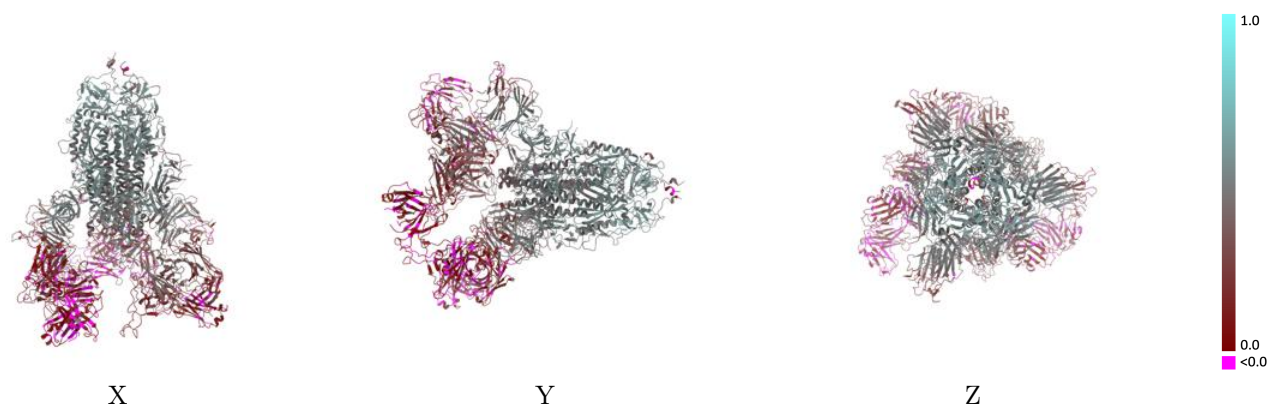
Y



Z

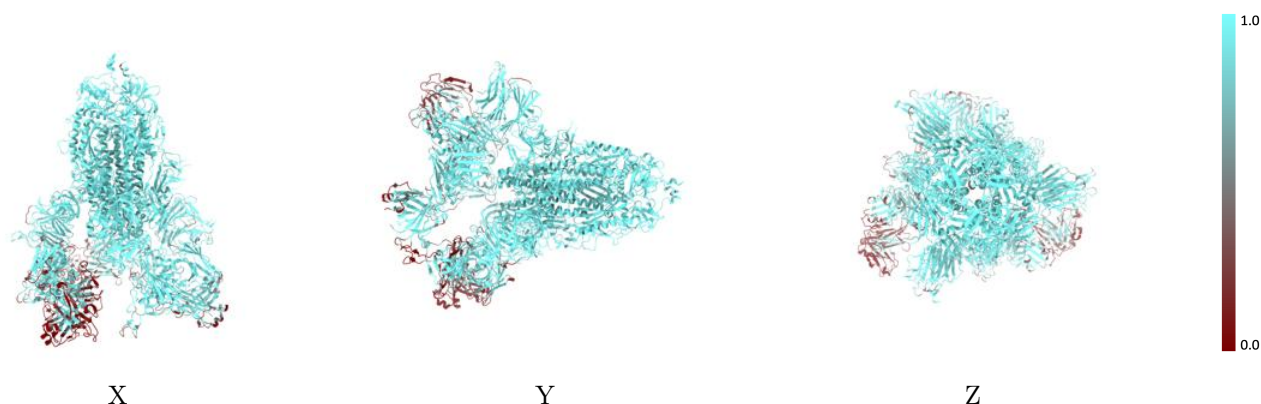
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



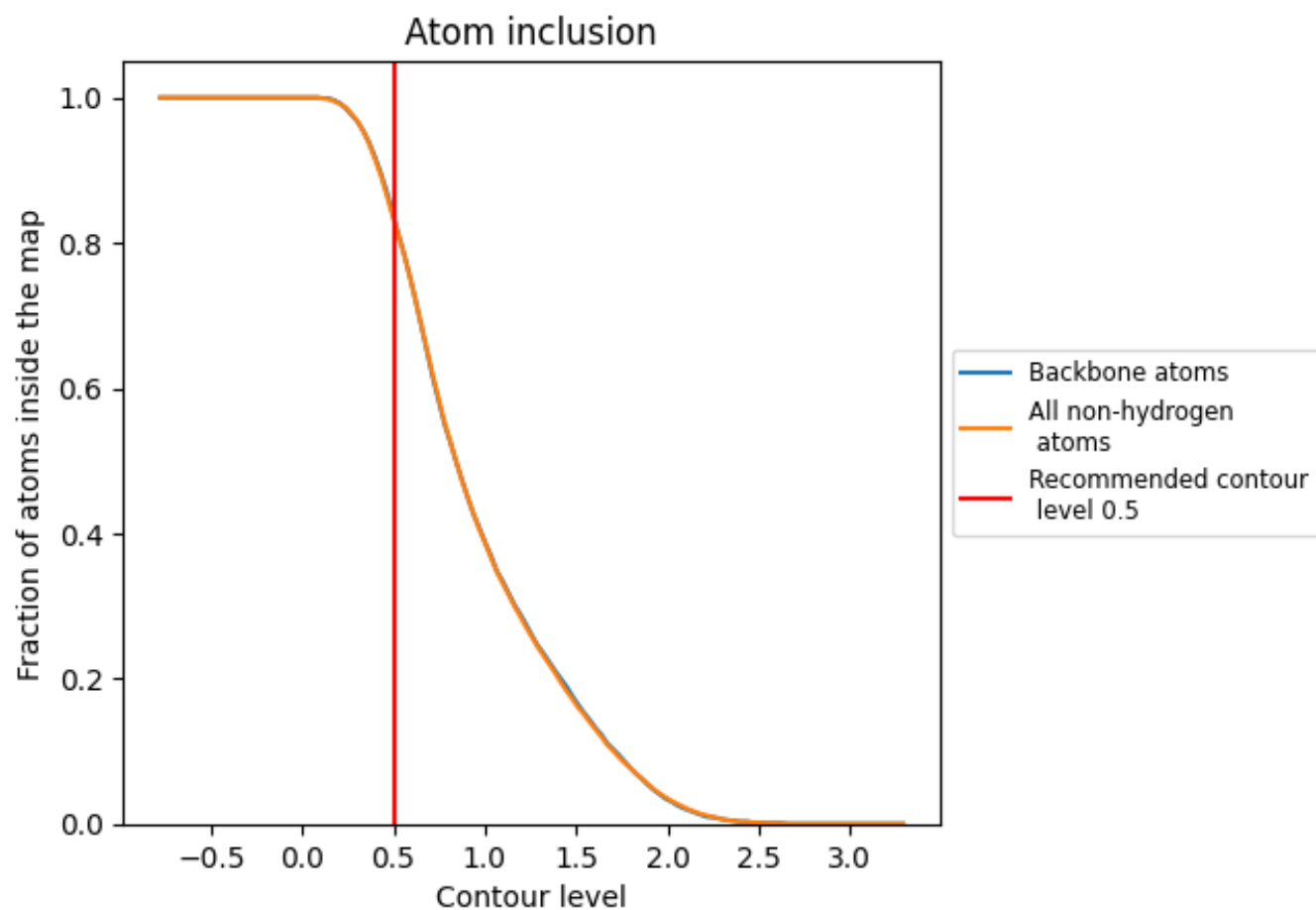
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 84% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8350	0.3330
A	0.9410	0.4140
B	0.9220	0.4280
C	0.9050	0.3910
H	0.8700	0.2130
I	0.5890	0.1110
J	0.4450	0.0740
L	0.8380	0.2130
M	0.3990	0.0480
N	0.5240	0.1120

1.0

0.0

<0.0