



wwPDB EM Validation Summary Report ⓘ

Mar 5, 2026 – 04:45 AM UTC

PDB ID : 7ZCF / pdb_00007zcf
EMDB ID : EMD-14629
Title : SARS-CoV-2 Spike RBD in complex with the single chain fragment scFv76
(Focused Refinement)
Authors : Chaves-Sanjuan, A.; Bolognesi, M.
Deposited on : 2022-03-28
Resolution : 4.00 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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

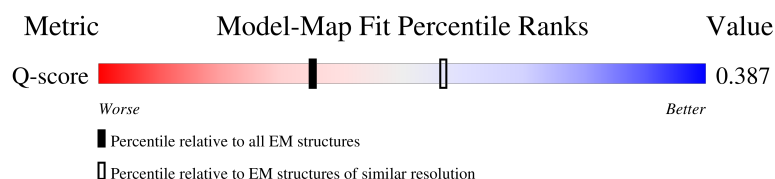
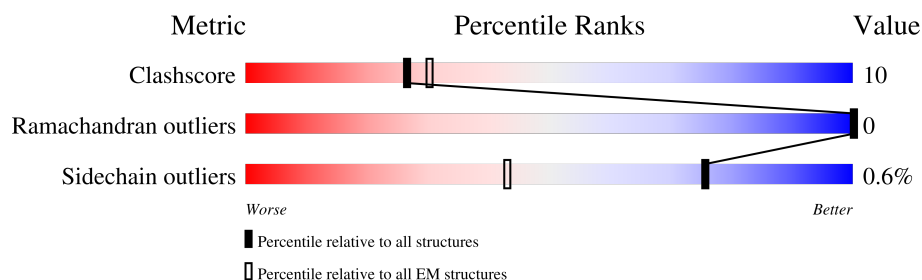
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.00 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	7587 (3.50 - 4.50)

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	C	1288	11% 84%
2	E	264	9% 67% 18% 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3312 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 Spike glycoprotein,Fibritin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	201	Total	C	N	O	S	0	0
			1595	1025	266	296	8		

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	682	GLY	ARG	engineered mutation	UNP P0DTC2
C	683	SER	ARG	engineered mutation	UNP P0DTC2
C	685	SER	ARG	engineered mutation	UNP P0DTC2
C	817	PRO	PHE	engineered mutation	UNP P0DTC2
C	892	PRO	ALA	engineered mutation	UNP P0DTC2
C	899	PRO	ALA	engineered mutation	UNP P0DTC2
C	942	PRO	ALA	engineered mutation	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	1209	GLY	-	linker	UNP P0DTC2
C	1210	SER	-	linker	UNP P0DTC2
C	1232	LEU	PHE	conflict	UNP P10104
C	1238	GLY	-	expression tag	UNP P10104
C	1239	ARG	-	expression tag	UNP P10104
C	1240	SER	-	expression tag	UNP P10104
C	1241	LEU	-	expression tag	UNP P10104
C	1242	GLU	-	expression tag	UNP P10104
C	1243	VAL	-	expression tag	UNP P10104
C	1244	LEU	-	expression tag	UNP P10104
C	1245	PHE	-	expression tag	UNP P10104
C	1246	GLN	-	expression tag	UNP P10104
C	1247	GLY	-	expression tag	UNP P10104
C	1248	PRO	-	expression tag	UNP P10104
C	1249	GLY	-	expression tag	UNP P10104
C	1250	HIS	-	expression tag	UNP P10104
C	1251	HIS	-	expression tag	UNP P10104
C	1252	HIS	-	expression tag	UNP P10104
C	1253	HIS	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1254	HIS	-	expression tag	UNP P10104
C	1255	HIS	-	expression tag	UNP P10104
C	1256	HIS	-	expression tag	UNP P10104
C	1257	HIS	-	expression tag	UNP P10104
C	1258	SER	-	expression tag	UNP P10104
C	1259	ALA	-	expression tag	UNP P10104
C	1260	TRP	-	expression tag	UNP P10104
C	1261	SER	-	expression tag	UNP P10104
C	1262	HIS	-	expression tag	UNP P10104
C	1263	PRO	-	expression tag	UNP P10104
C	1264	GLN	-	expression tag	UNP P10104
C	1265	PHE	-	expression tag	UNP P10104
C	1266	GLU	-	expression tag	UNP P10104
C	1267	LYS	-	expression tag	UNP P10104
C	1268	GLY	-	expression tag	UNP P10104
C	1269	GLY	-	expression tag	UNP P10104
C	1270	GLY	-	expression tag	UNP P10104
C	1271	SER	-	expression tag	UNP P10104
C	1272	GLY	-	expression tag	UNP P10104
C	1273	GLY	-	expression tag	UNP P10104
C	1274	GLY	-	expression tag	UNP P10104
C	1275	GLY	-	expression tag	UNP P10104
C	1276	SER	-	expression tag	UNP P10104
C	1277	GLY	-	expression tag	UNP P10104
C	1278	GLY	-	expression tag	UNP P10104
C	1279	SER	-	expression tag	UNP P10104
C	1280	ALA	-	expression tag	UNP P10104
C	1281	TRP	-	expression tag	UNP P10104
C	1282	SER	-	expression tag	UNP P10104
C	1283	HIS	-	expression tag	UNP P10104
C	1284	PRO	-	expression tag	UNP P10104
C	1285	GLN	-	expression tag	UNP P10104
C	1286	PHE	-	expression tag	UNP P10104
C	1287	GLU	-	expression tag	UNP P10104
C	1288	LYS	-	expression tag	UNP P10104

- Molecule 2 is a protein called scFv76 single chain fragment.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	225	Total	C	N	O	S	0	0
			1703	1074	289	334	6		

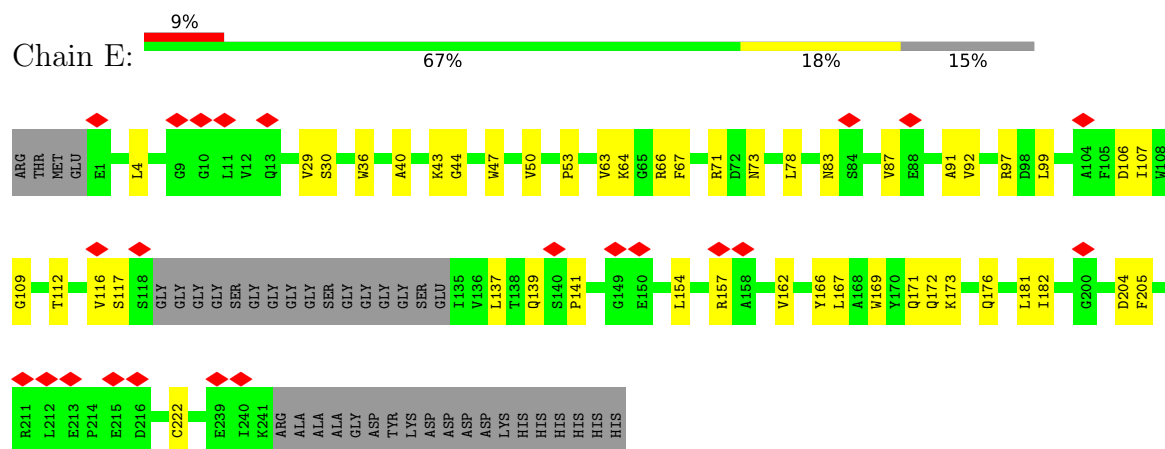
- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
3	C	1	14	8	1	5	0

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

- Molecule 2: scFv76 single chain fragment



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	87623	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	120000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	4.014	Depositor
Minimum map value	-3.897	Depositor
Average map value	-0.004	Depositor
Map value standard deviation	0.044	Depositor
Recommended contour level	0.6	Depositor
Map size ($\text{\AA}$)	355.6, 355.6, 355.6	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.889, 0.889, 0.889	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	C	0.12	0/1641	0.38	0/2233
2	E	0.15	0/1740	0.39	0/2362
All	All	0.14	0/3381	0.39	0/4595

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

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	C	1595	0	1518	36	0
2	E	1703	0	1661	30	0
3	C	14	0	13	0	0
All	All	3312	0	3192	65	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 10.

The worst 5 of 65 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:C:354:ASN:OD1	1:C:399:SER:OG	2.09	0.70
1:C:331:ASN:OD1	1:C:332:ILE:N	2.23	0.67
1:C:353:TRP:HE1	1:C:466:ARG:HB3	1.60	0.67
2:E:47:TRP:HE1	2:E:50:VAL:HG23	1.60	0.67
1:C:472:ILE:HG22	1:C:474:GLN:H	1.62	0.64

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	C	199/1288 (16%)	166 (83%)	33 (17%)	0	100	100
2	E	221/264 (84%)	192 (87%)	29 (13%)	0	100	100
All	All	420/1552 (27%)	358 (85%)	62 (15%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	174/1116 (16%)	174 (100%)	0	100	100
2	E	184/207 (89%)	182 (99%)	2 (1%)	65	74
All	All	358/1323 (27%)	356 (99%)	2 (1%)	76	81

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	73	ASN
2	E	139	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	E	83	ASN
2	E	160	GLN
2	E	224	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	C	1301	1	14,14,15	0.26	0	17,19,21	0.55	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1301	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1301	NAG	O5-C5-C6-O6
3	C	1301	NAG	C4-C5-C6-O6
3	C	1301	NAG	C1-C2-N2-C7
3	C	1301	NAG	C3-C2-N2-C7

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.

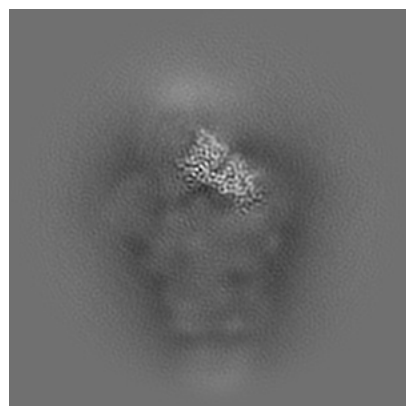
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14629. These allow visual inspection of the internal detail of the map and identification of artifacts.

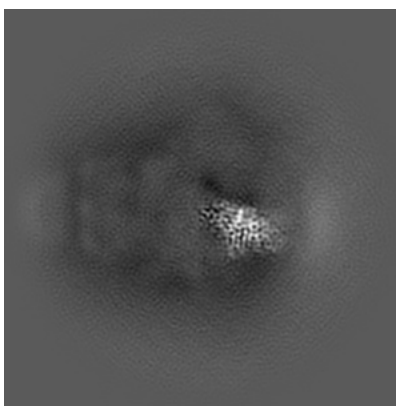
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

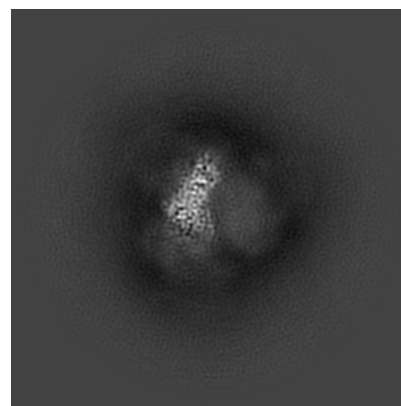
6.1.1 Primary map



X

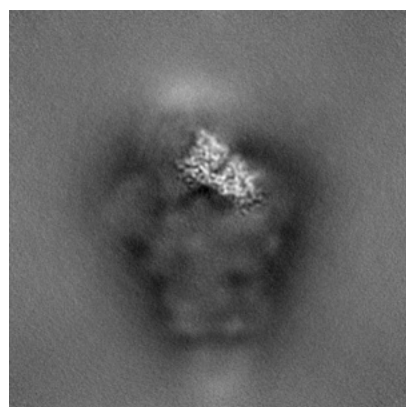


Y

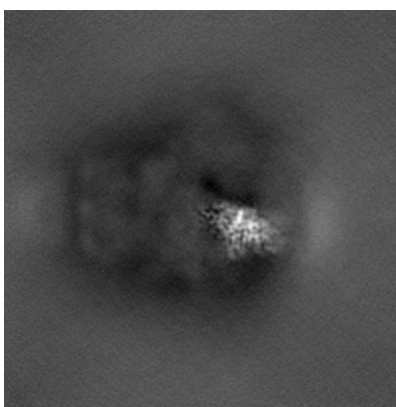


Z

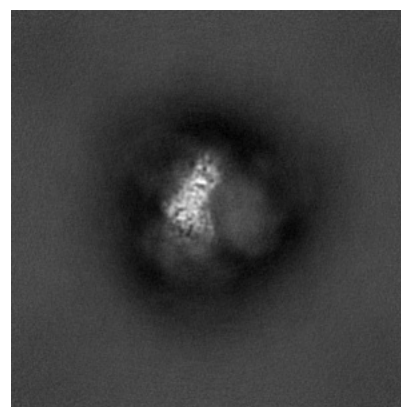
6.1.2 Raw map



X



Y

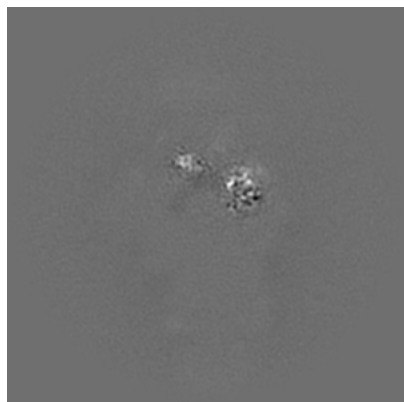


Z

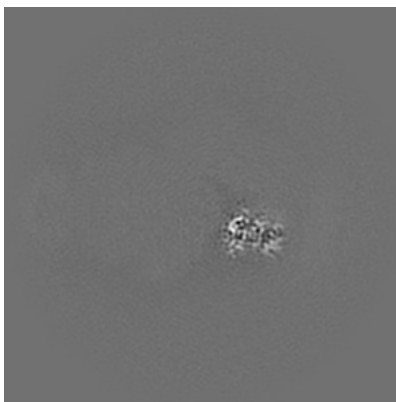
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

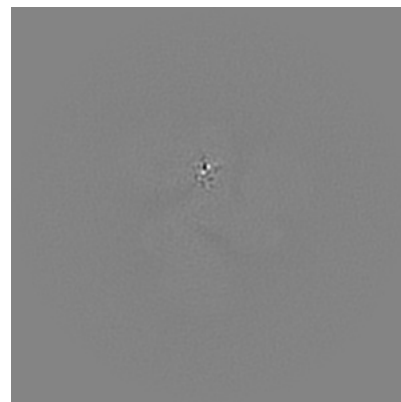
6.2.1 Primary map



X Index: 200

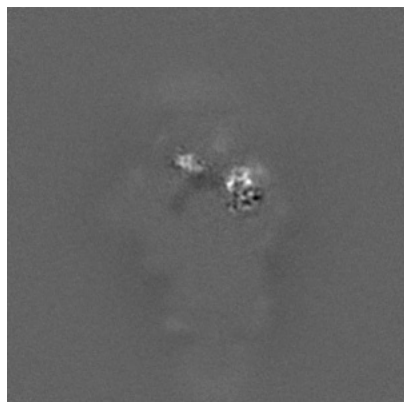


Y Index: 200

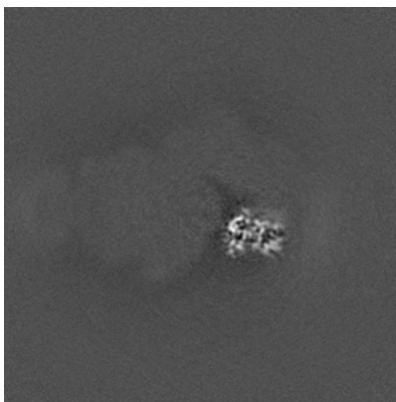


Z Index: 200

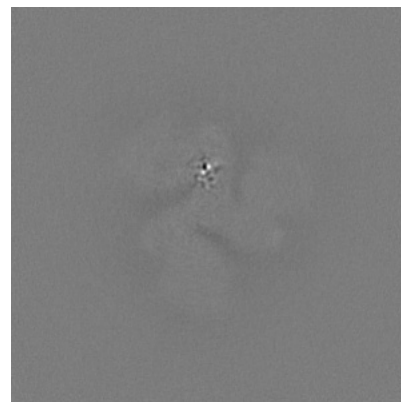
6.2.2 Raw map



X Index: 200



Y Index: 200

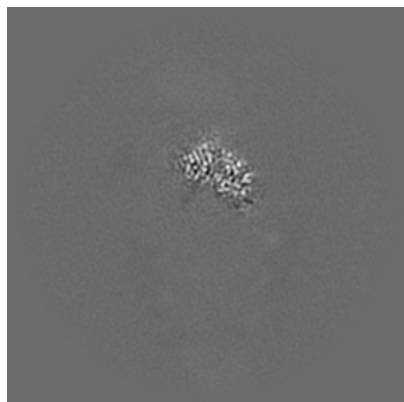


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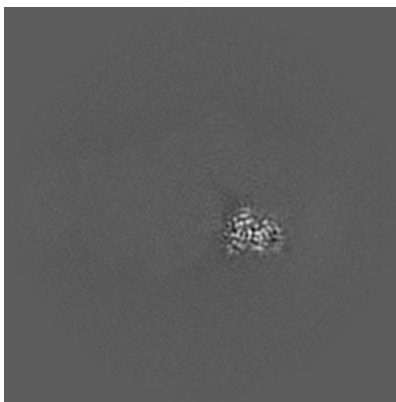
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

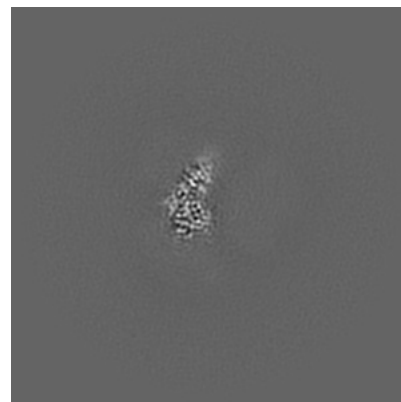
6.3.1 Primary map



X Index: 184

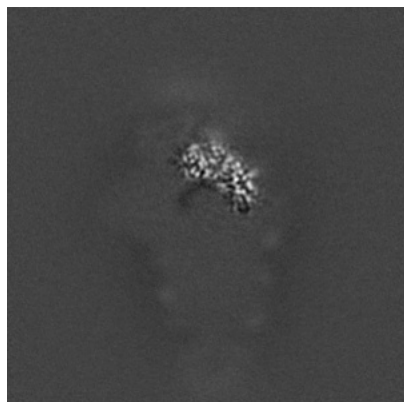


Y Index: 197

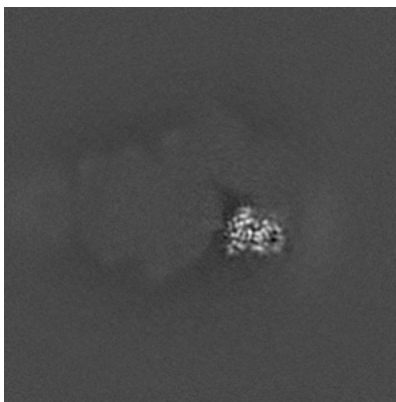


Z Index: 235

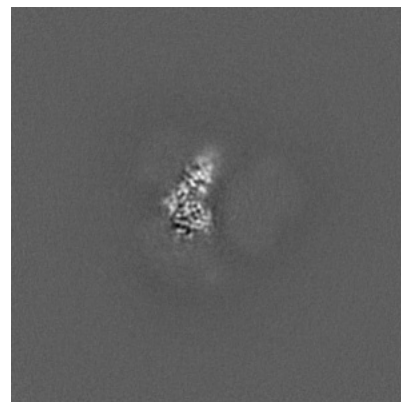
6.3.2 Raw map



X Index: 187



Y Index: 197

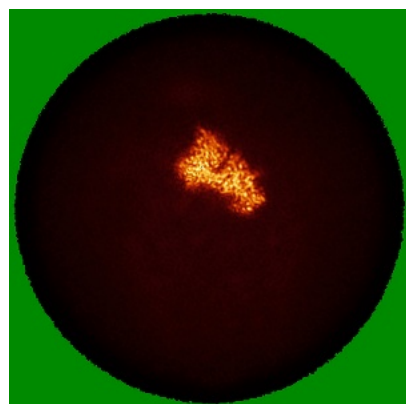


Z Index: 235

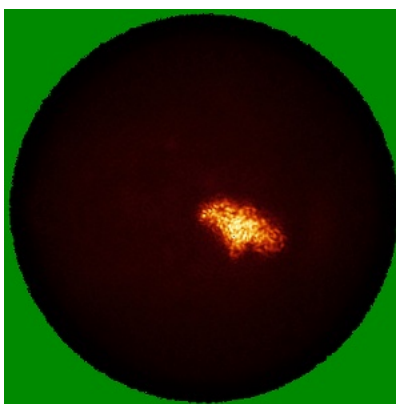
The images above show the largest variance slices of the map in three orthogonal directions.

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

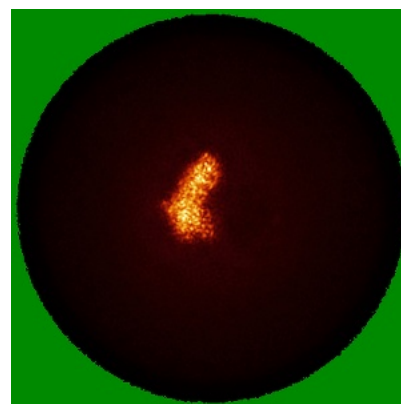
6.4.1 Primary map



X

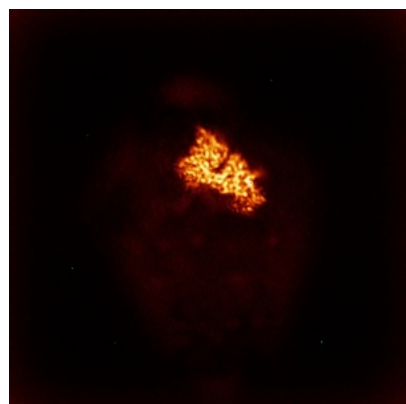


Y

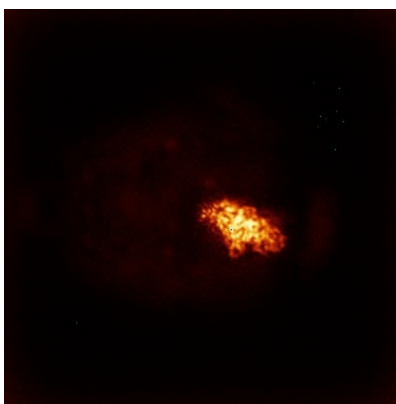


Z

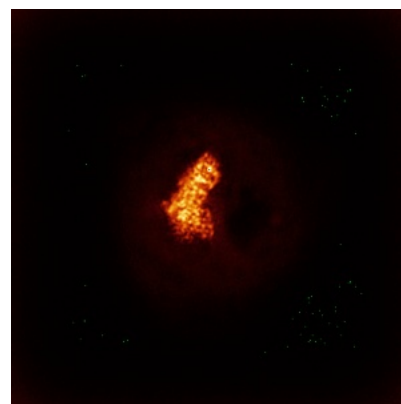
6.4.2 Raw map



X



Y

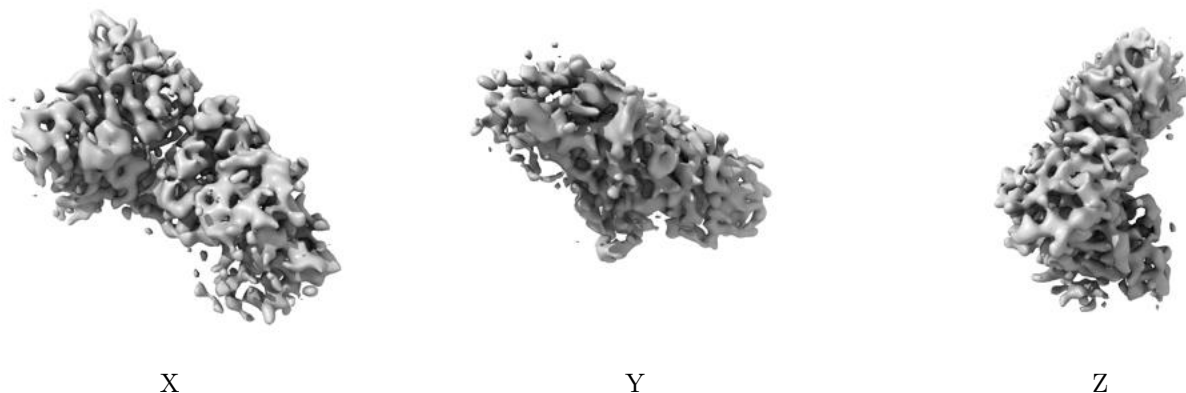


Z

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.

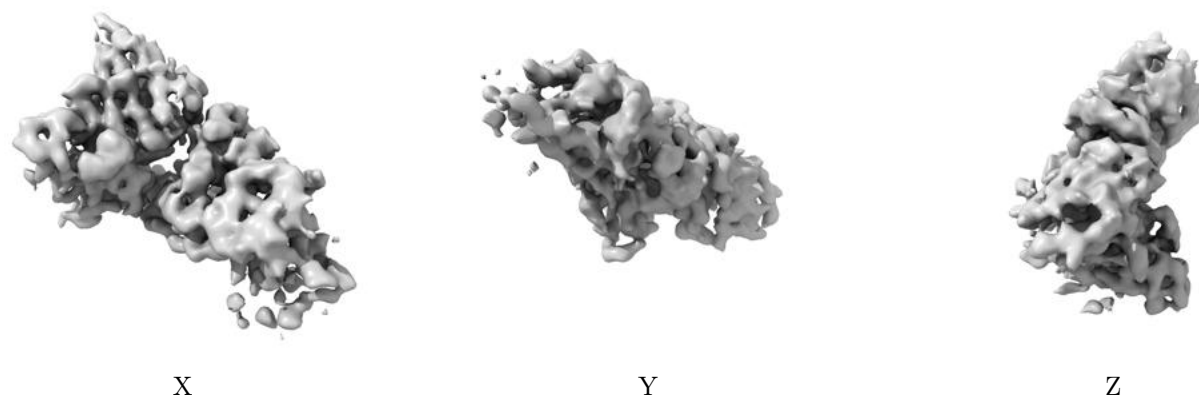
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.6. 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.

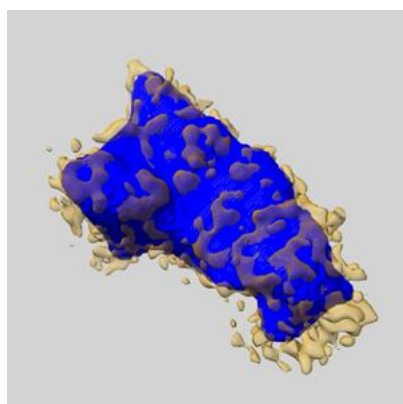
6.6 Mask visualisation [i](#)

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

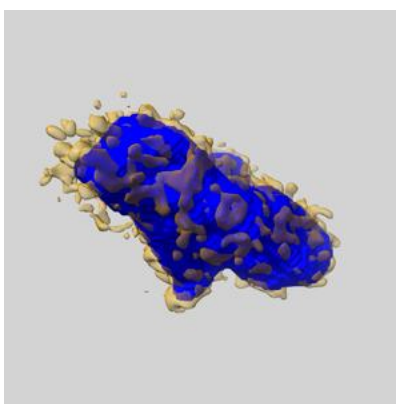
A mask typically either:

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

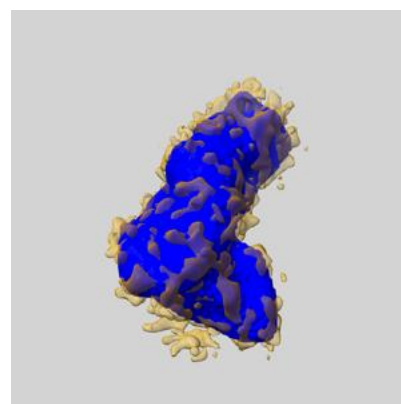
6.6.1 emd_14629_msk_1.map [i](#)



X



Y

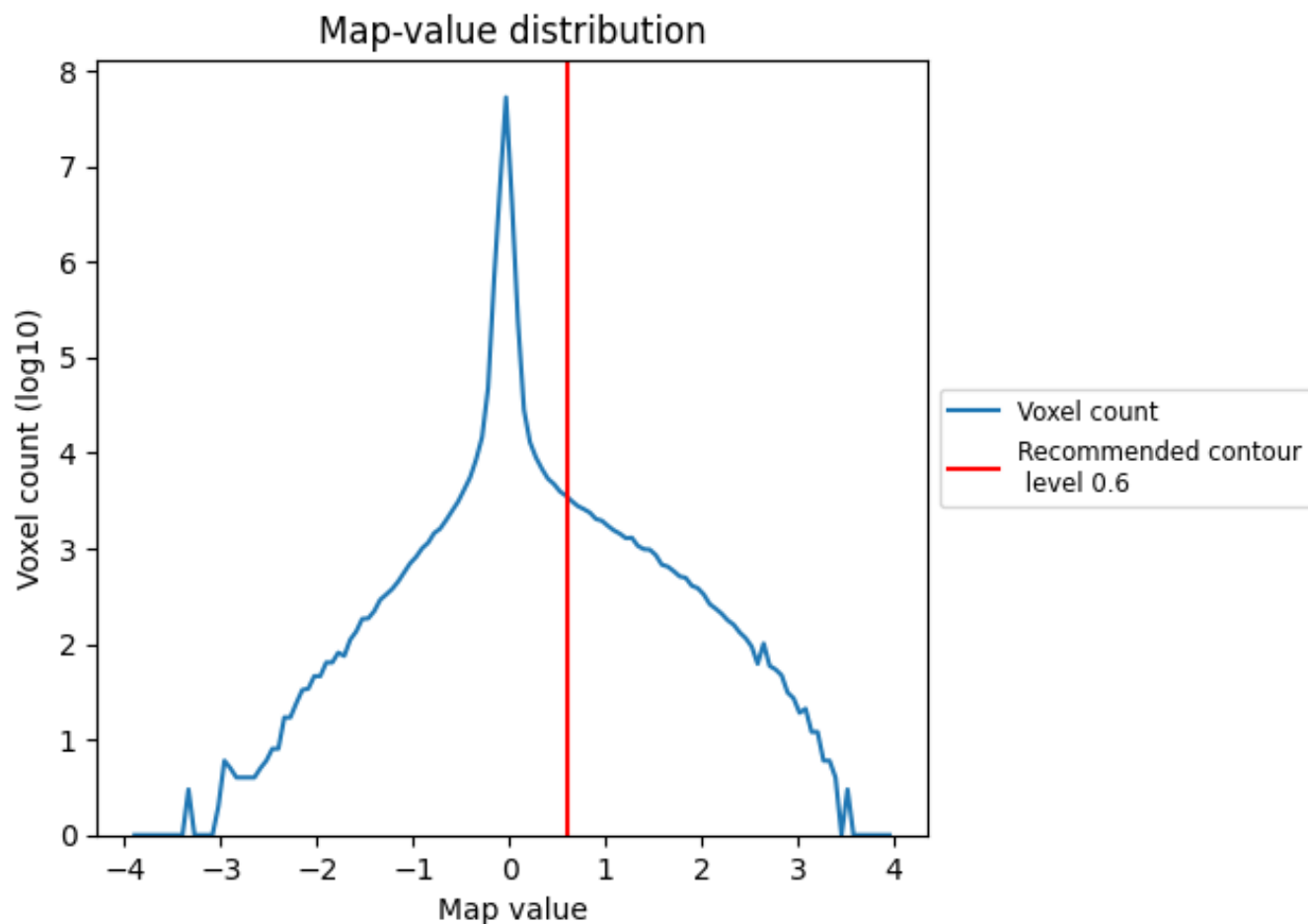


Z

7 Map analysis [i](#)

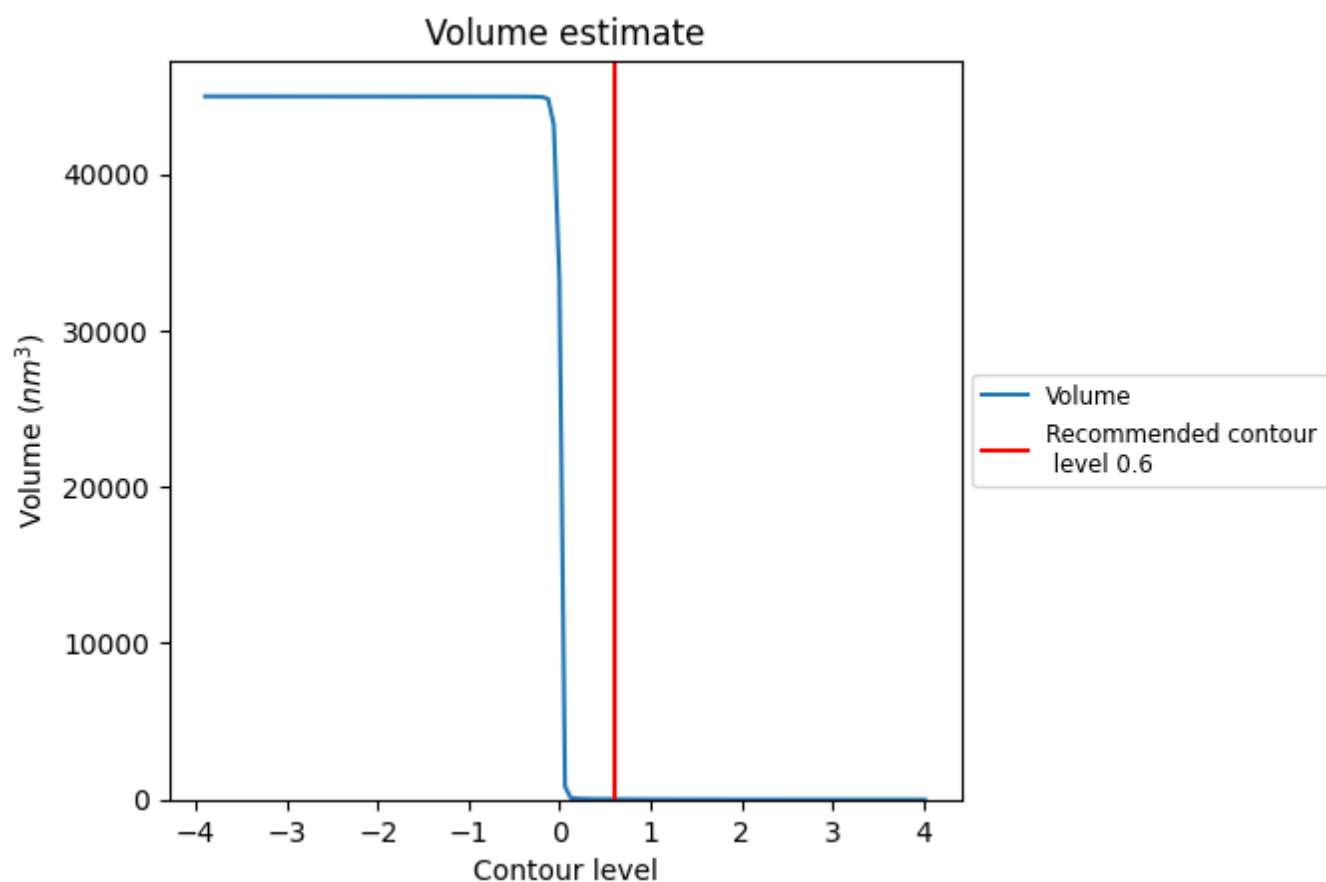
This section contains the results of statistical analysis of the map.

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.

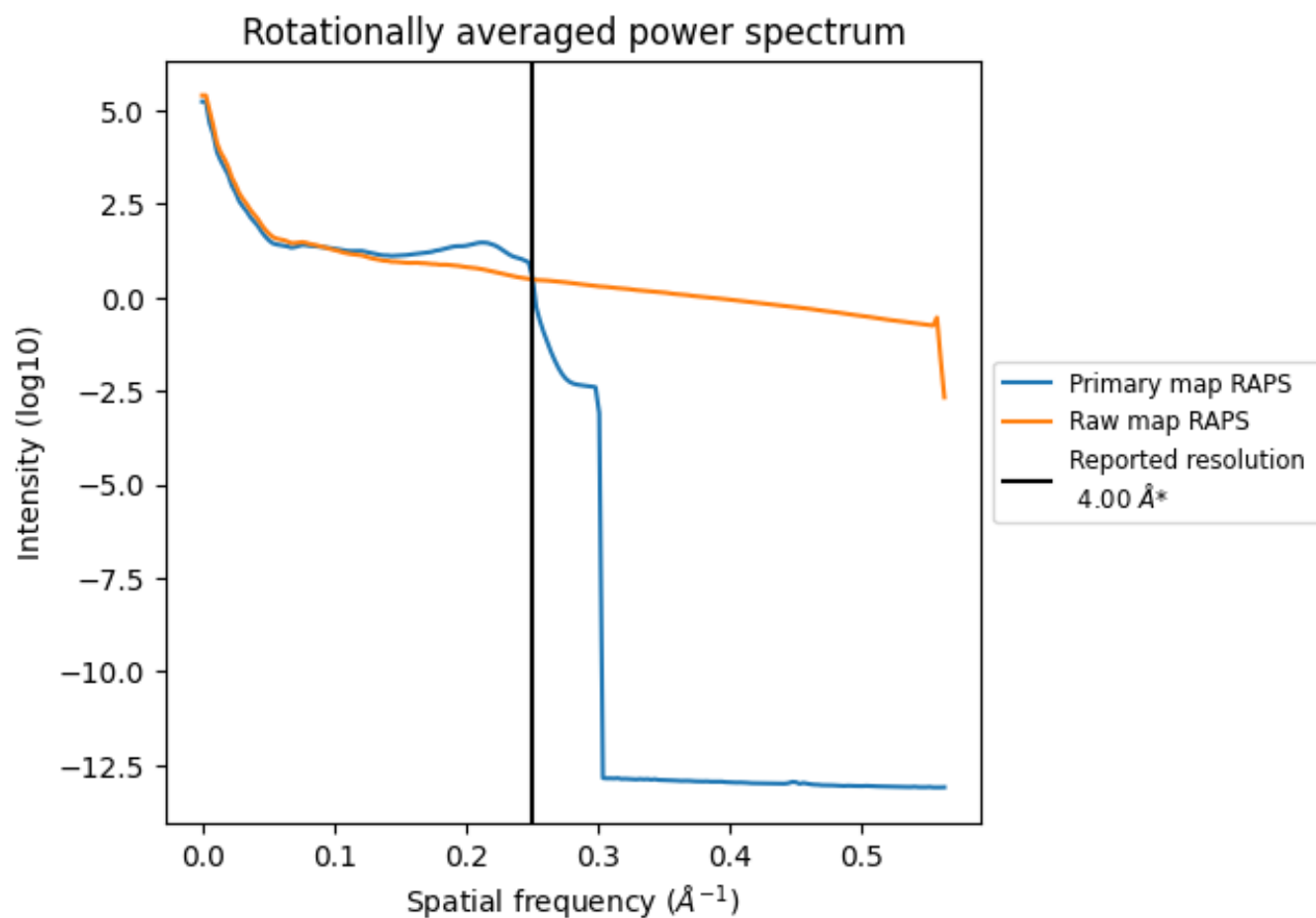
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 24 nm³; this corresponds to an approximate mass of 22 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](#)

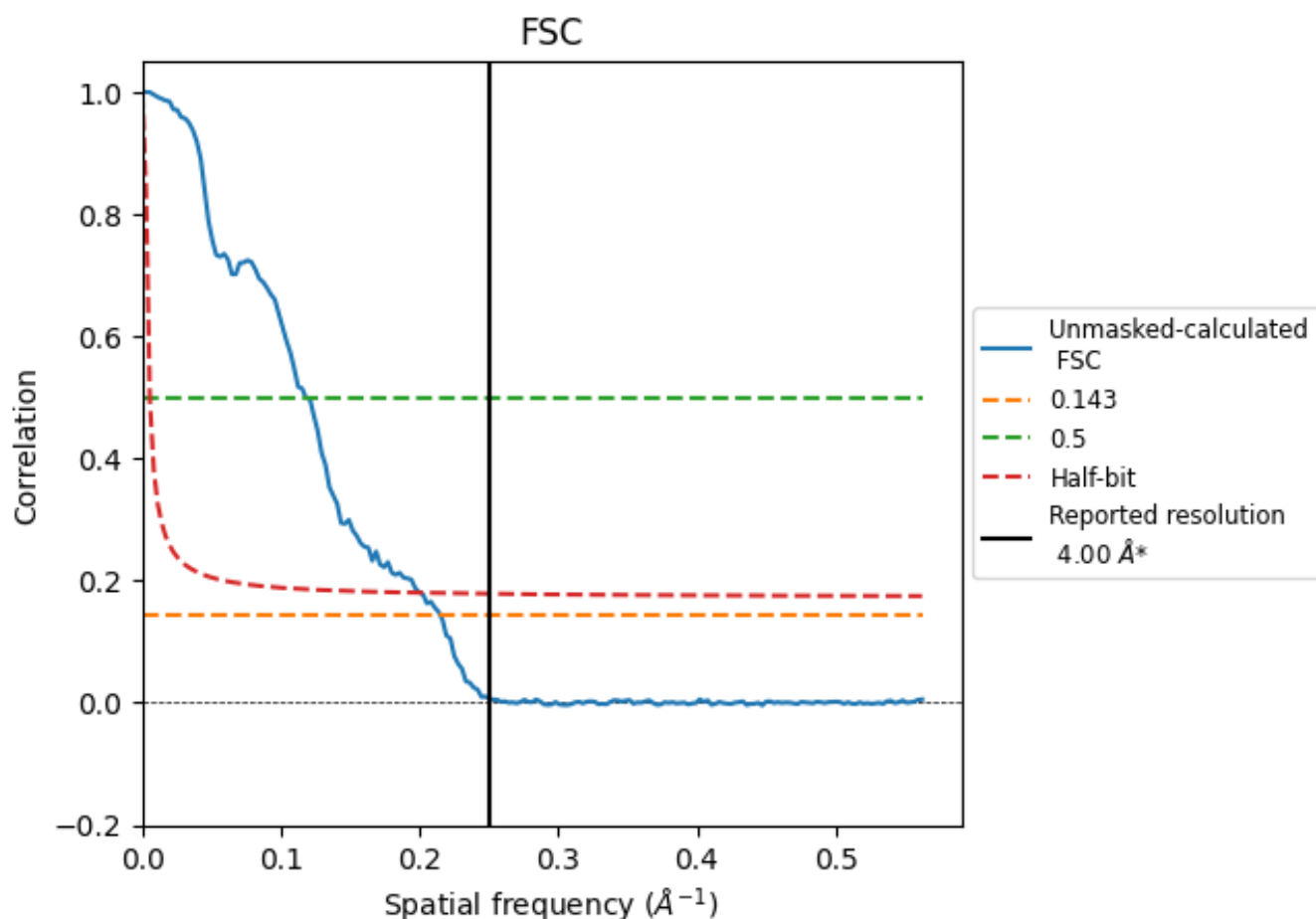


*Reported resolution corresponds to spatial frequency of $0.250 \text{ \AA}^{-1}$

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 0.250 Å⁻¹

8.2 Resolution estimates [i](#)

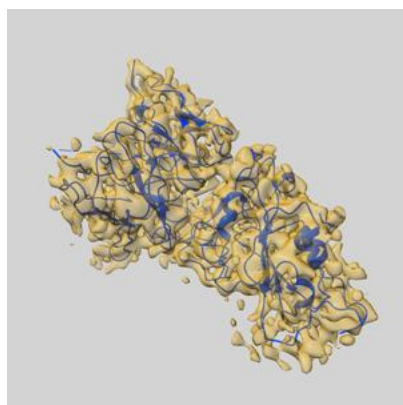
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.00	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.65	8.49	4.99

*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 4.65 differs from the reported value 4.0 by more than 10 %

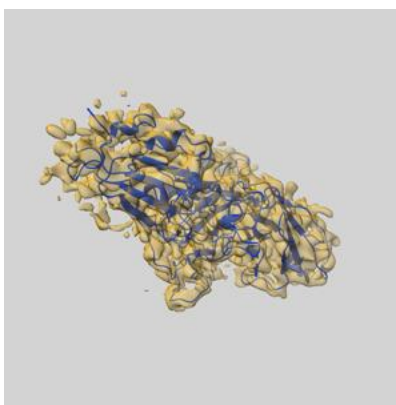
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-14629 and PDB model 7ZCF. Per-residue inclusion information can be found in section [3](#) on page [6](#).

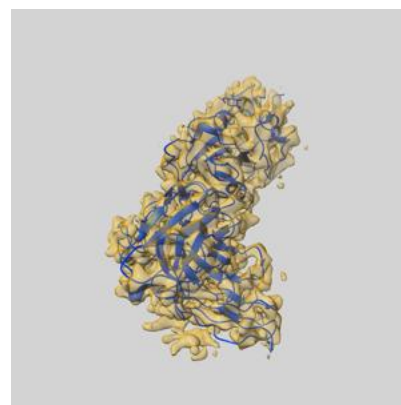
9.1 Map-model overlay [i](#)



X



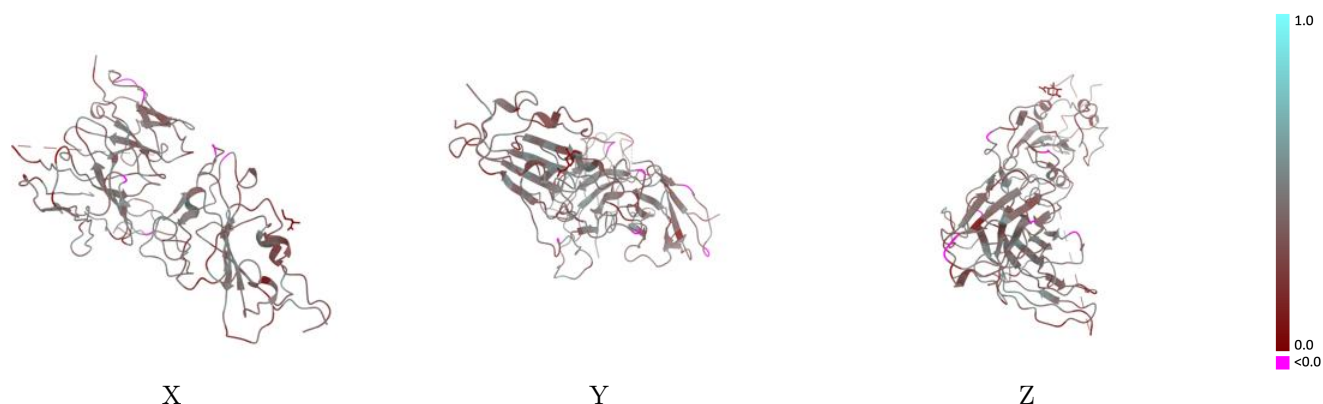
Y



Z

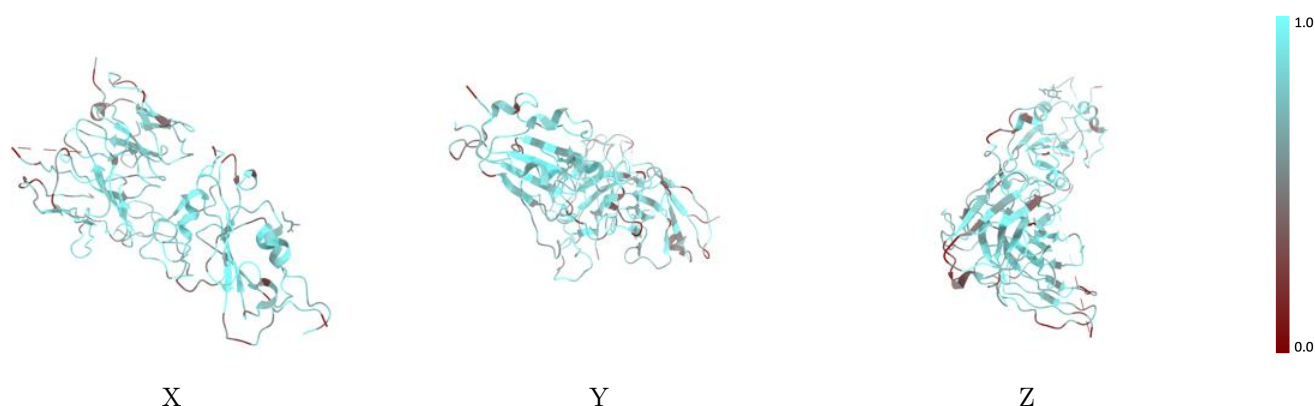
The images above show the 3D surface view of the map at the recommended contour level 0.6 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



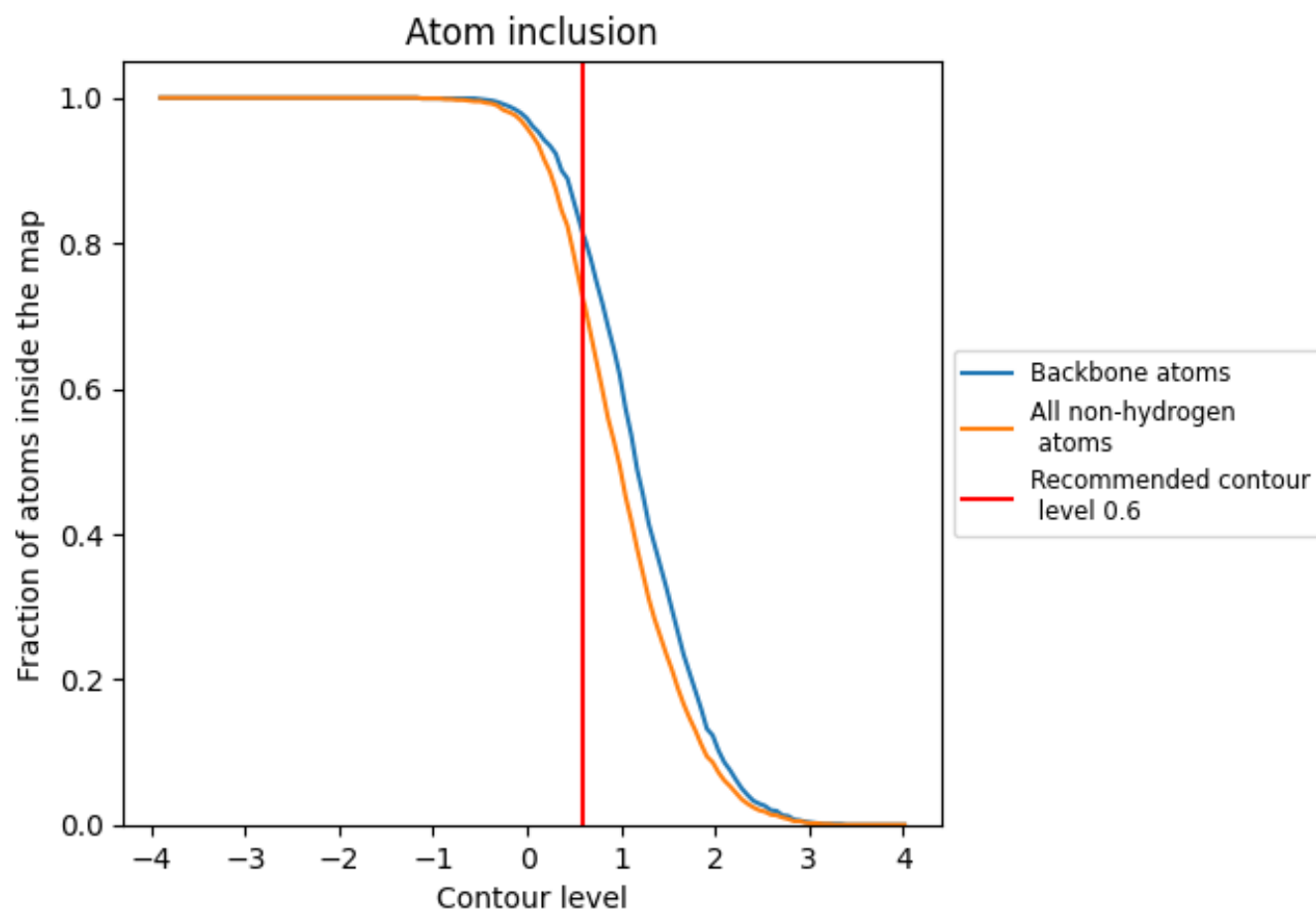
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.6).

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 72% 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 (0.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7200	0.3870
C	0.7290	0.3840
E	0.7120	0.3900

