wwPDB EM Validation Summary Report

Oct 18, 2026 – 05:37 PM JST

PDB ID	:	21RW / pdb_000021rw
EMDB ID	:	EMD-67953
Title	:	cryoEM structure of PROTEIN TANC1
Authors	:	Xue, C.Y.; Liu, Z.M.; Liu, W.; Zhong, D.
Deposited on	:	2025-12-23
Resolution	:	3.40 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:

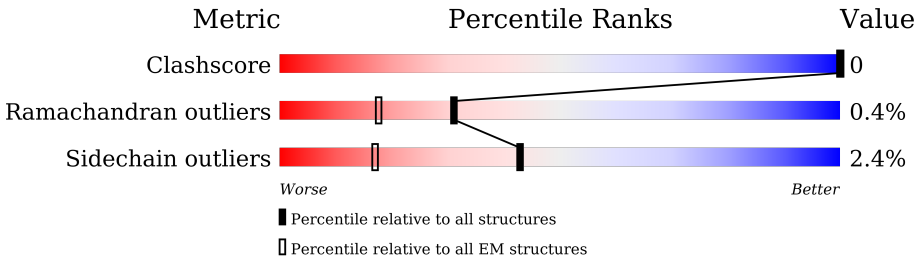
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

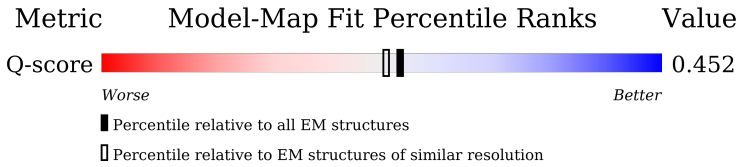
1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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	14717 (2.90 - 3.90)

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.

Mol	Chain	Length	Quality of chain
1	A	1856	38% 60%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 11627 atoms, of which 5851 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 Protein TANC1.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	738	11627	3673	5851	1023	1046	34	0	0

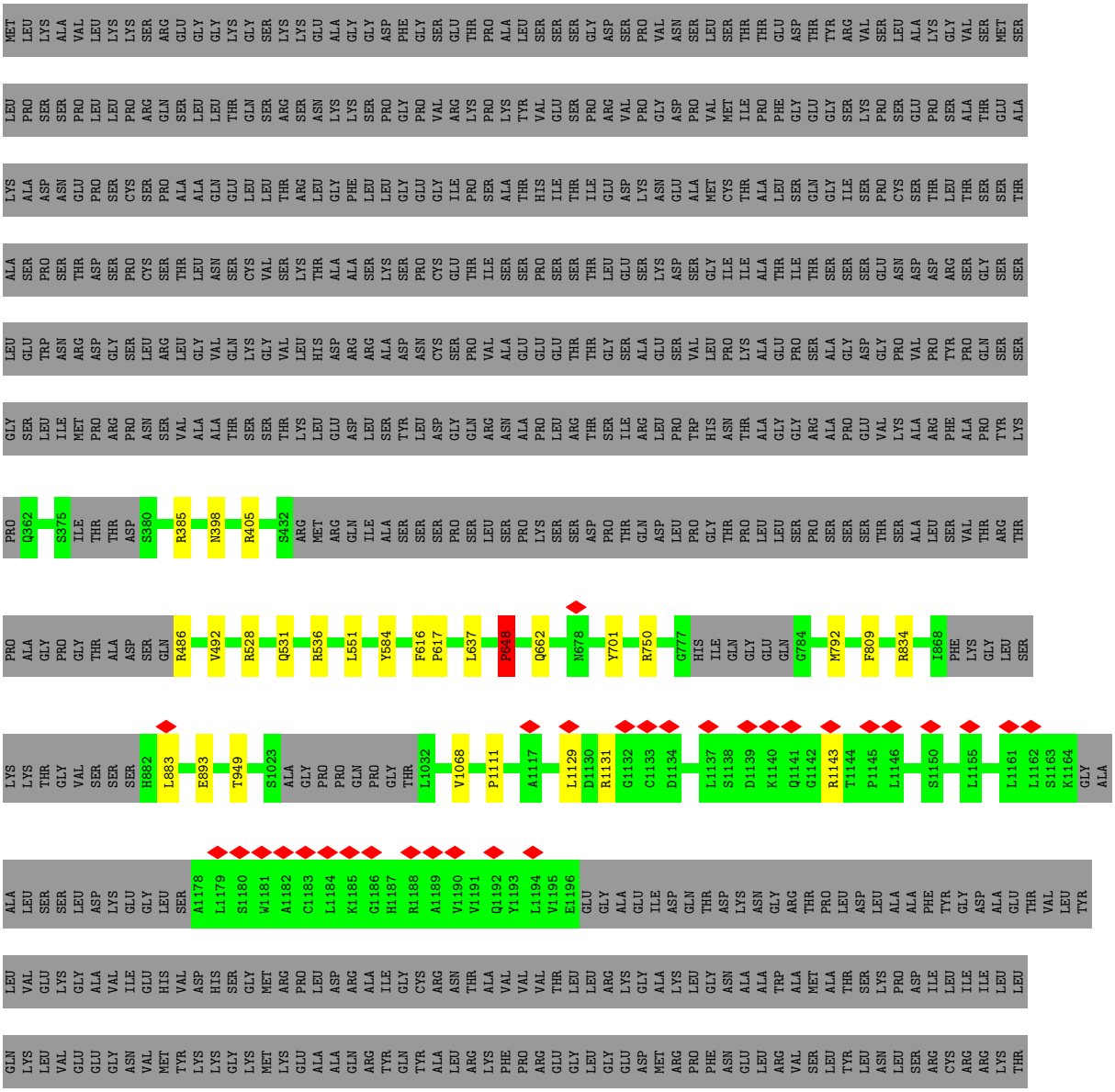
3 Residue-property plots (i)

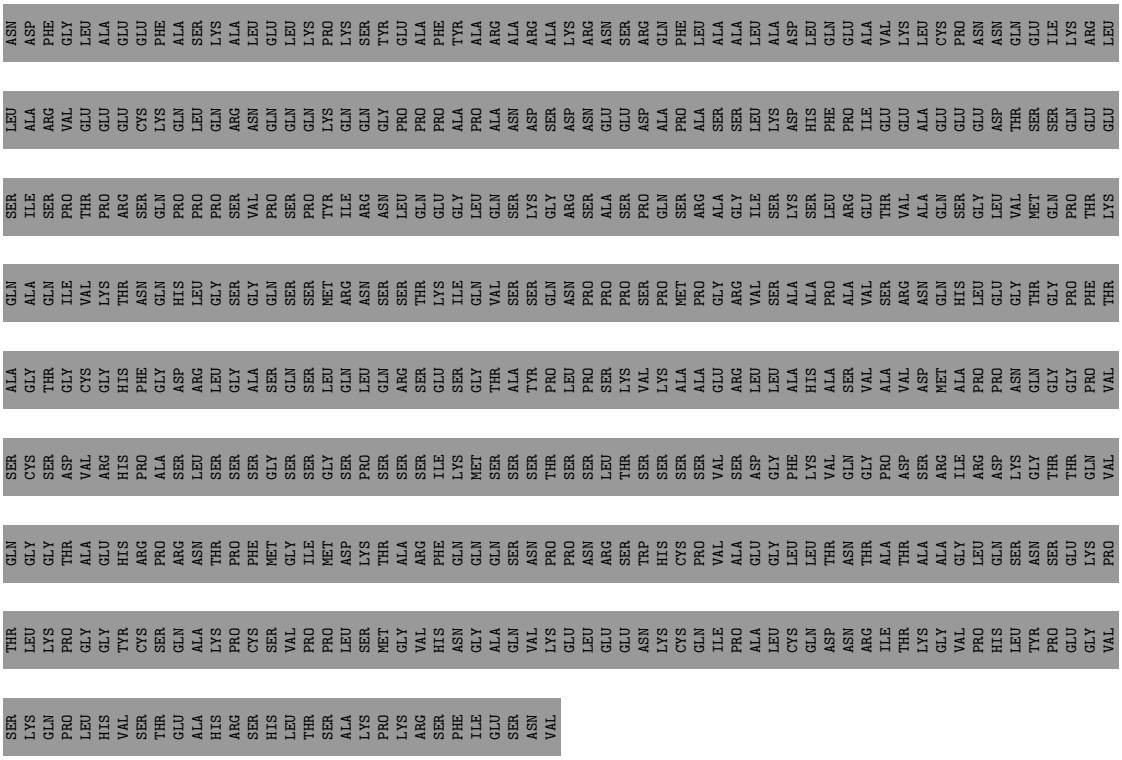
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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.

- Molecule 1: Protein TANC1

Chain A:






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

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, POINT, POINT, POINT	Depositor
Number of particles used	53000, 53000, 53000, 53000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION, PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI MORGAGNI	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.374	Depositor
Minimum map value	-0.239	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.0298	Depositor
Map size (Å)	189.36002, 189.36002, 189.36002	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.526, 0.526, 0.526	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	A	0.77	0/5886	1.33	12/7962 (0.2%)

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	A	0	2

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	648	PRO	CA-N-CD	-8.22	100.49	112.00
1	A	637	LEU	CA-C-N	5.59	126.83	119.84
1	A	637	LEU	C-N-CA	5.59	126.83	119.84
1	A	528	ARG	NE-CZ-NH2	5.56	124.20	119.20
1	A	834	ARG	NE-CZ-NH2	5.44	124.10	119.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	584	TYR	Sidechain
1	A	701	TYR	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	A	5776	5851	5844	0	0
All	All	5776	5851	5844	0	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.

There are no clashes within the asymmetric unit.

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	724/1856 (39%)	691 (95%)	30 (4%)	3 (0%)	30 59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	648	PRO
1	A	1068	VAL
1	A	617	PRO

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	A	628/1566 (40%)	613 (98%)	15 (2%)	43 62

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	662	GLN
1	A	949	THR
1	A	750	ARG
1	A	1129	LEU
1	A	883	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	977	HIS
1	A	1010	HIS
1	A	1064	HIS
1	A	763	HIS
1	A	769	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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

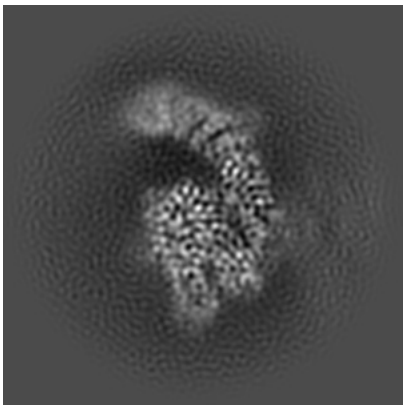
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-67953. These allow visual inspection of the internal detail of the map and identification of artifacts.

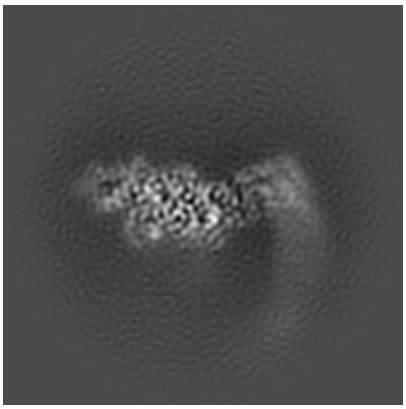
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

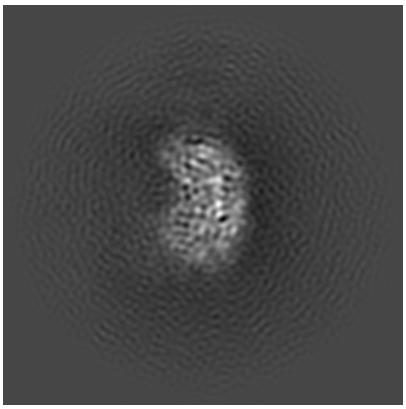
6.1.1 Primary map



X

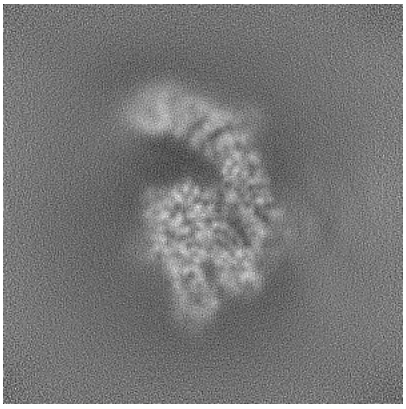


Y

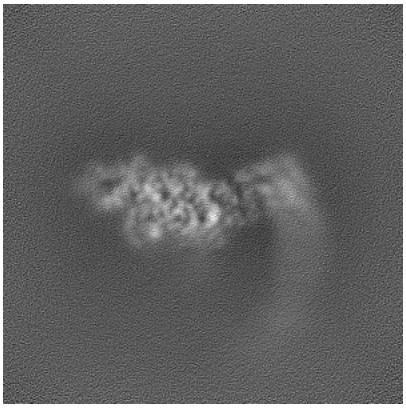


Z

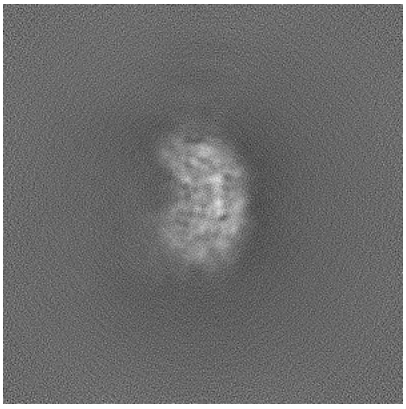
6.1.2 Raw map



X



Y

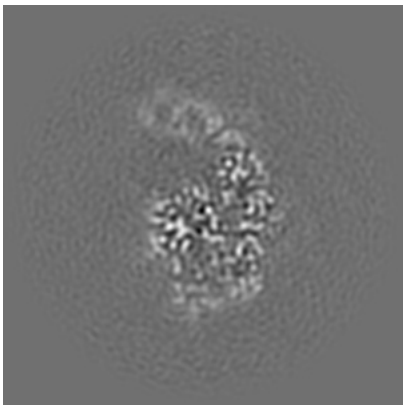


Z

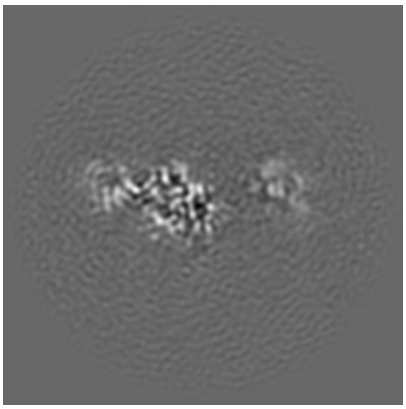
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

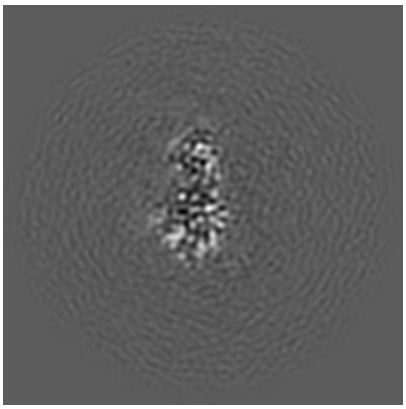
6.2.1 Primary map



X Index: 180

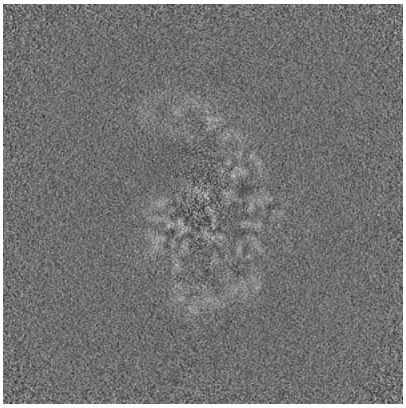


Y Index: 180

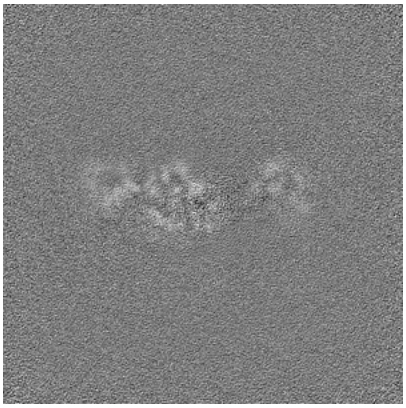


Z Index: 180

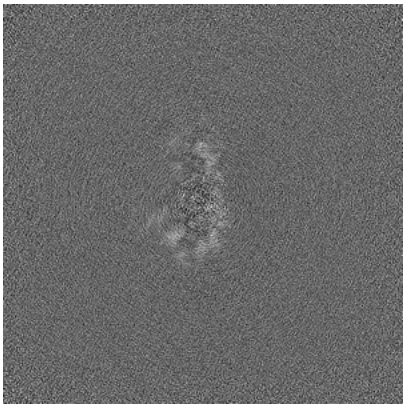
6.2.2 Raw map



X Index: 180



Y Index: 180

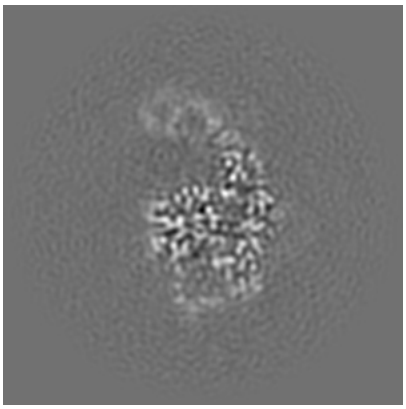


Z Index: 180

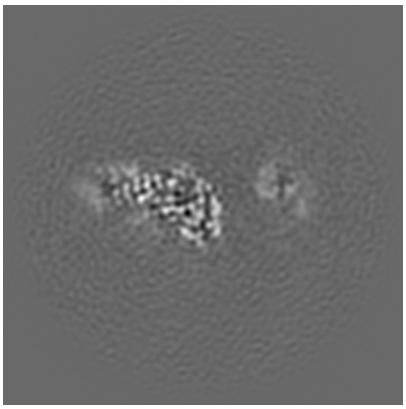
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

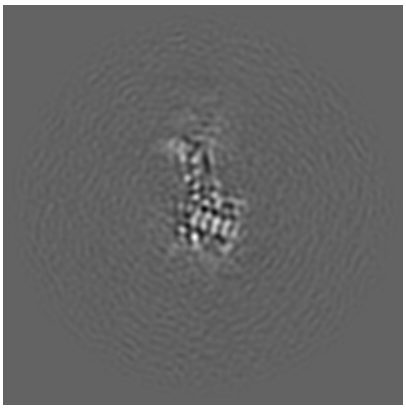
6.3.1 Primary map



X Index: 178

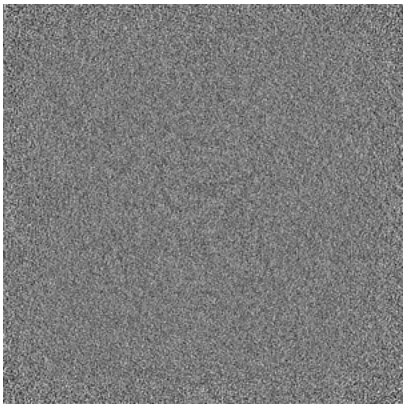


Y Index: 174

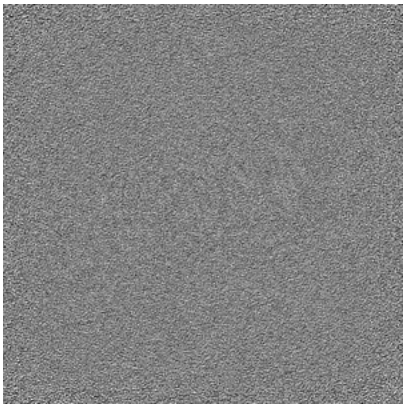


Z Index: 156

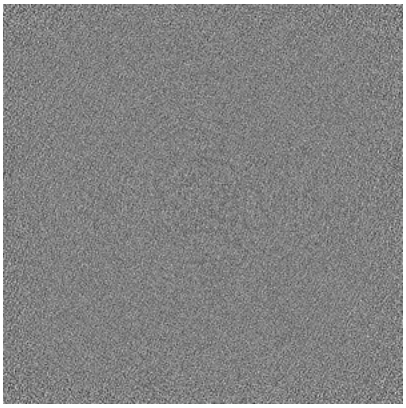
6.3.2 Raw map



X Index: 0



Y Index: 0

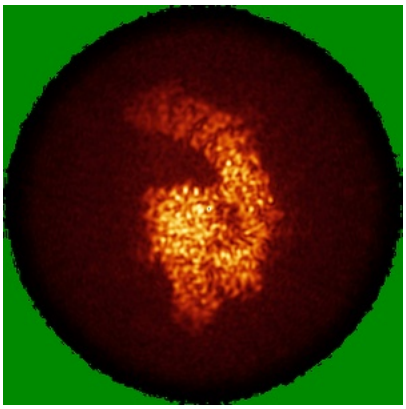


Z Index: 0

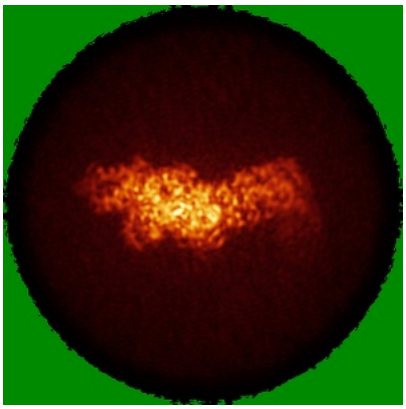
The images above show the largest variance slices of the map in three orthogonal directions.

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

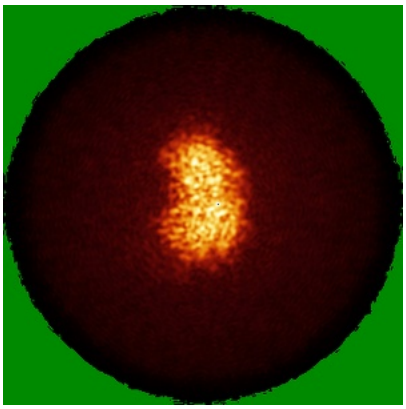
6.4.1 Primary map



X

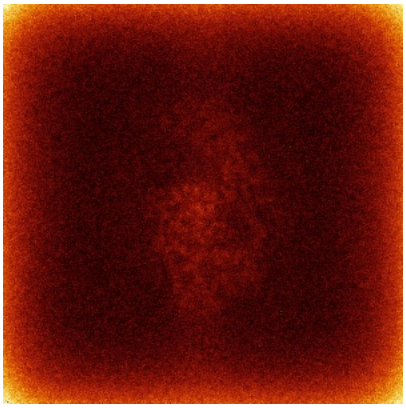


Y

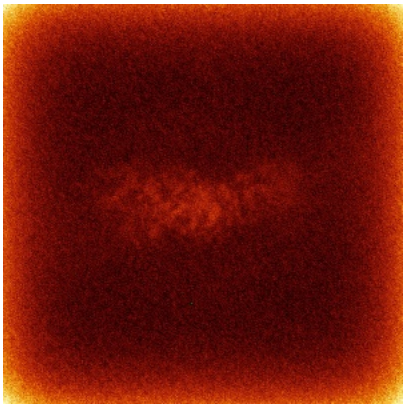


Z

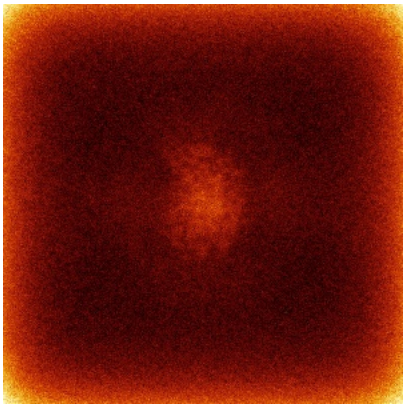
6.4.2 Raw map



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)

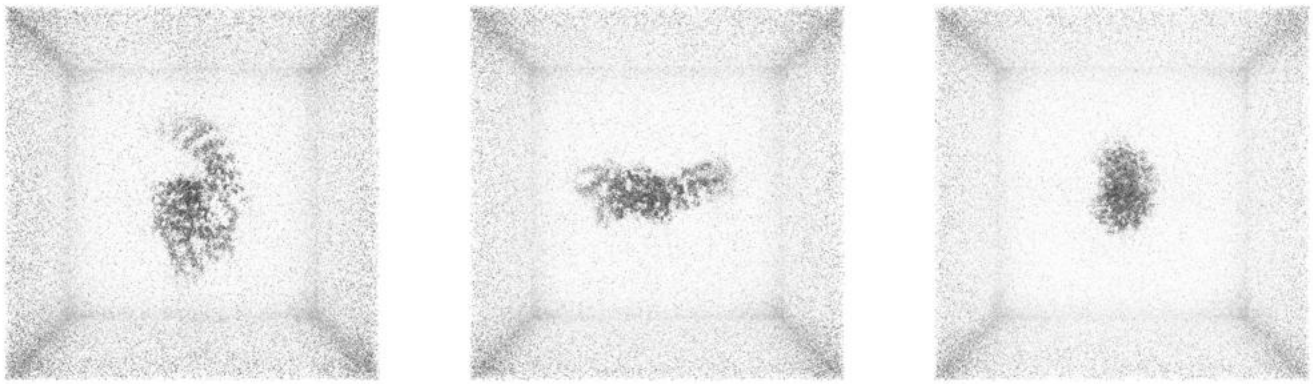
6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.0298. 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



X Y Z

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.

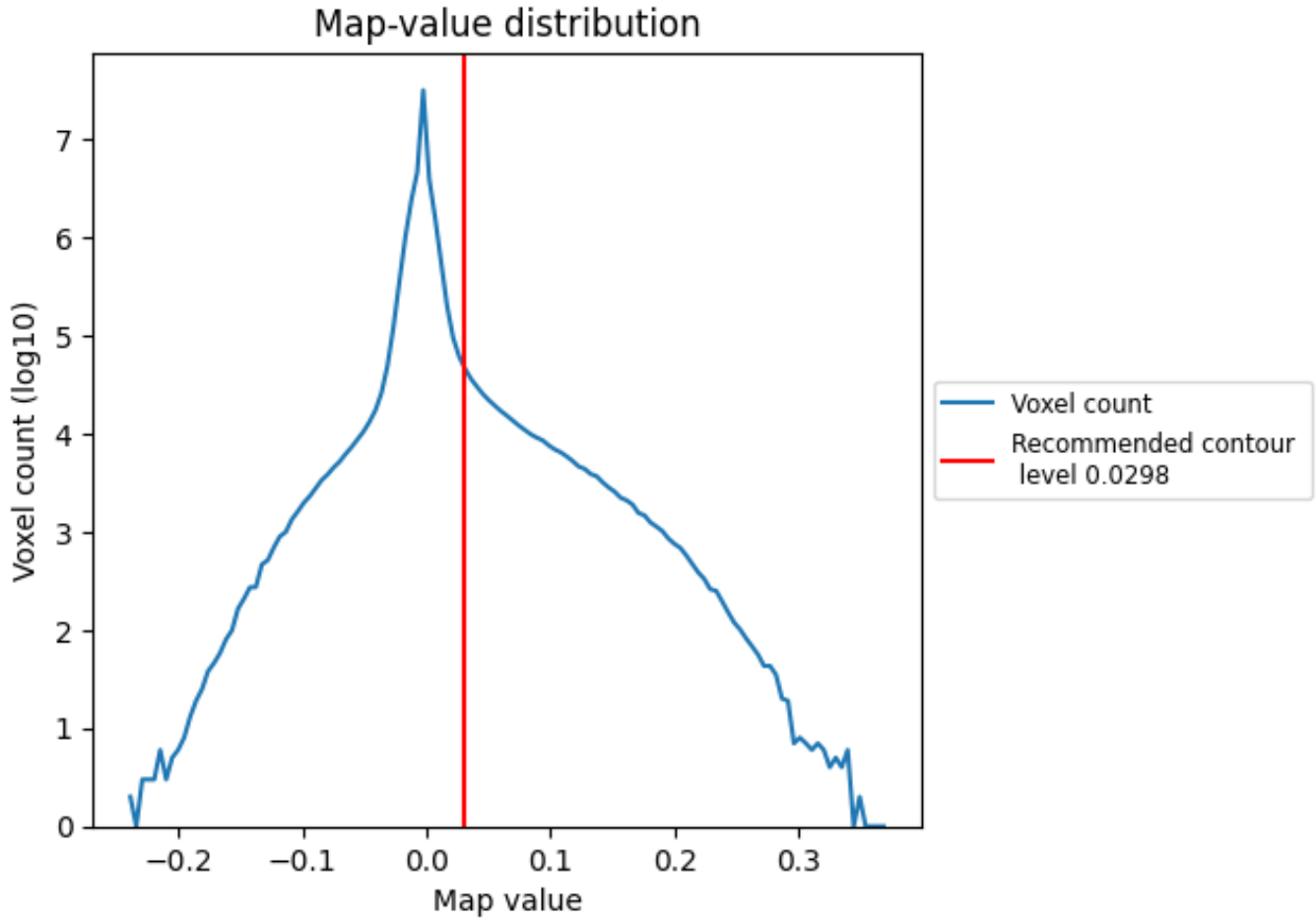
6.6 Mask visualisation (i)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis

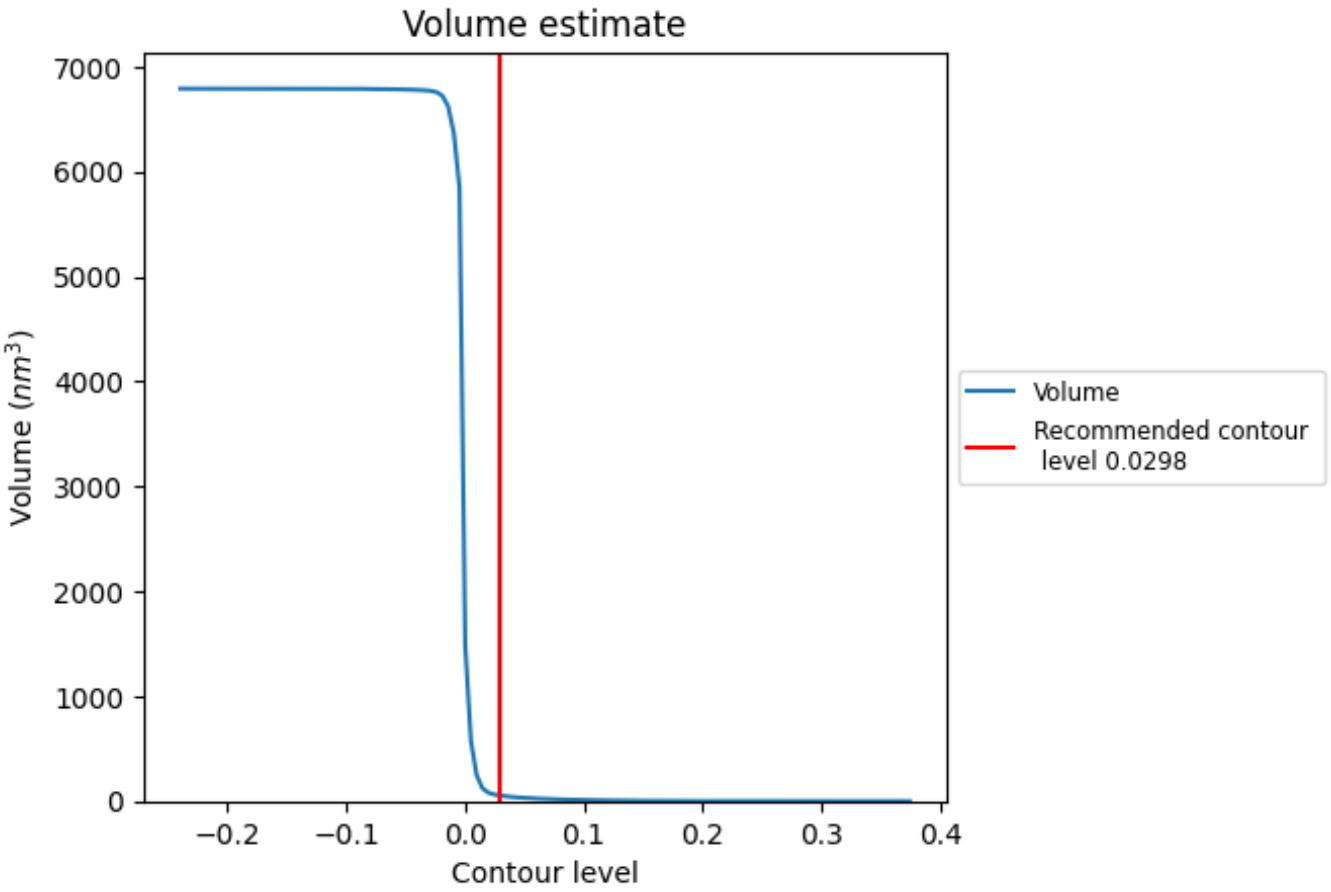
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

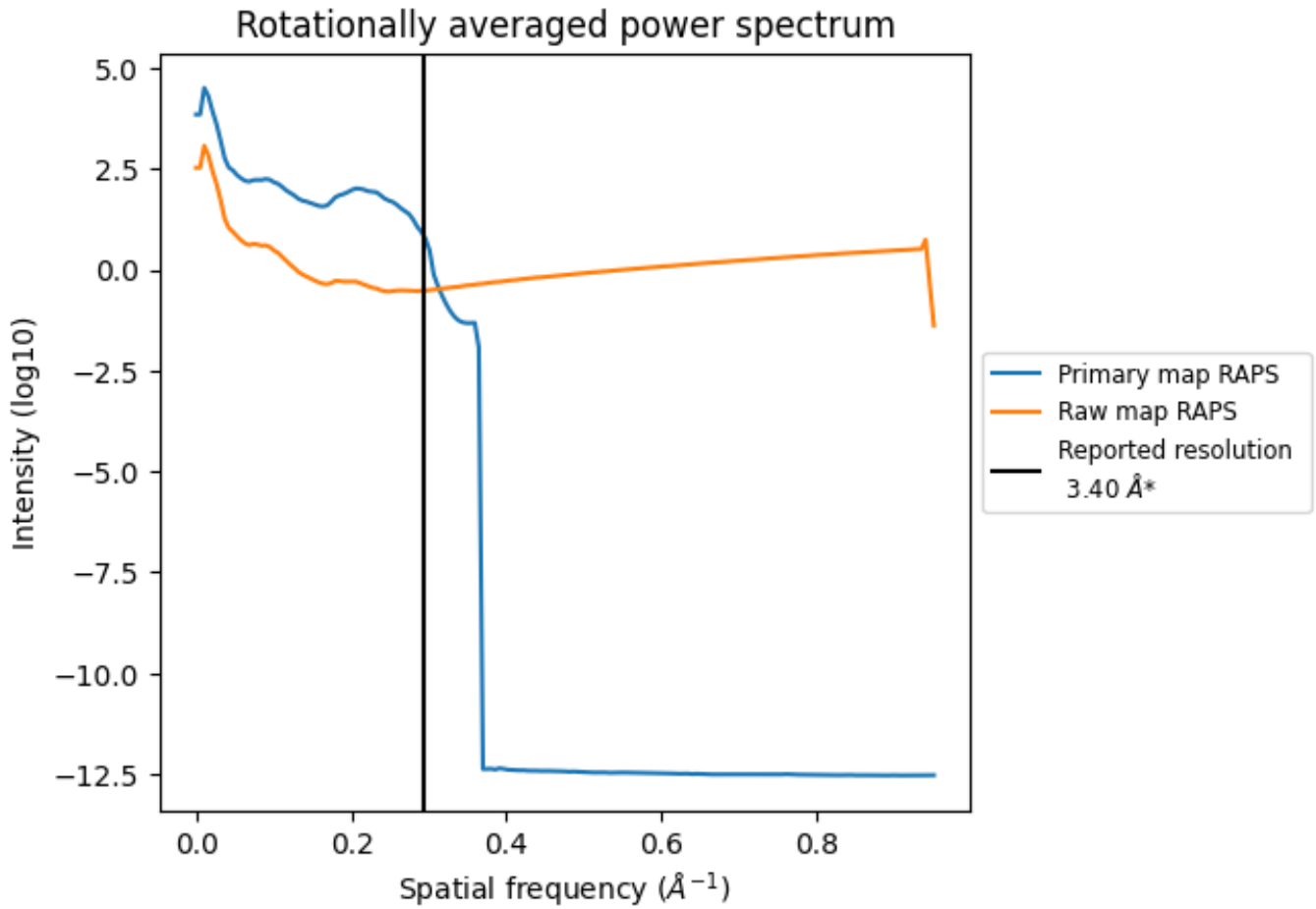
7.2 Volume estimate (i)



The volume at the recommended contour level is 53 nm^3; this corresponds to an approximate mass of 48 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)

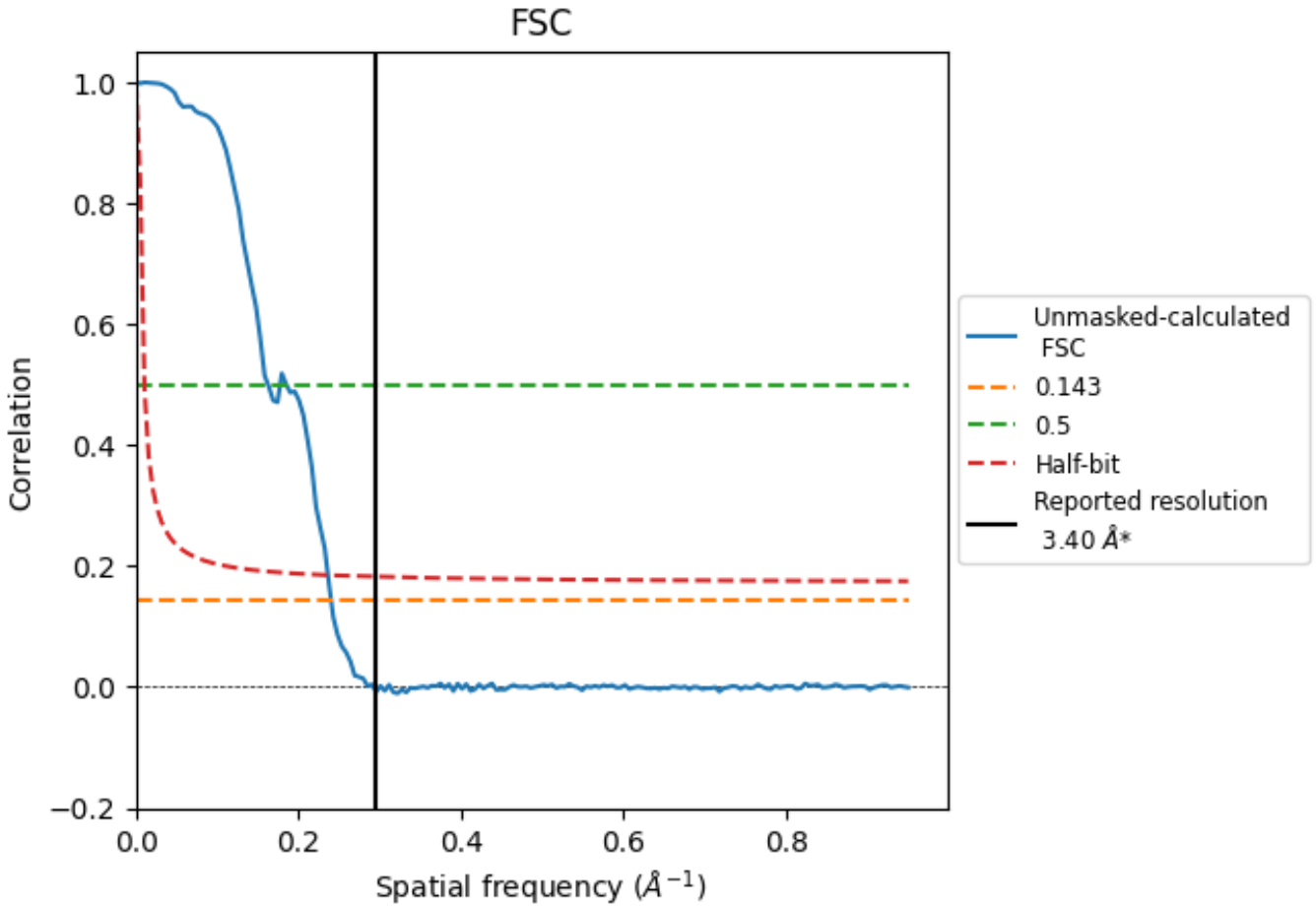


*Reported resolution corresponds to spatial frequency of 0.294 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.294 Å^{-1}

8.2 Resolution estimates (i)

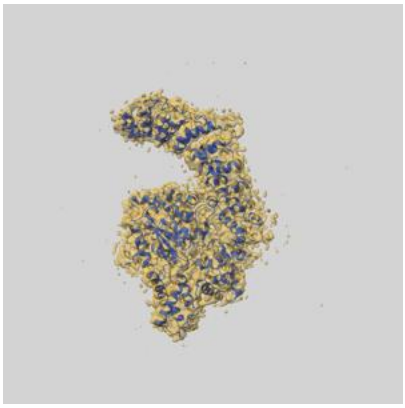
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	-	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.16	6.15	4.23

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

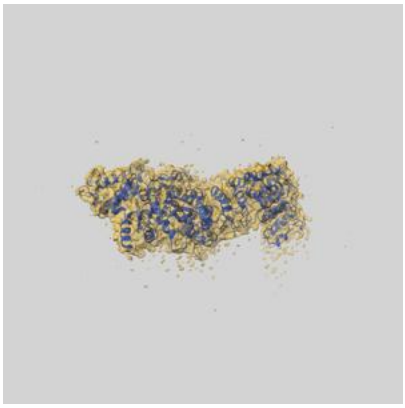
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-67953 and PDB model 21RW. Per-residue inclusion information can be found in section 3 on page 4.

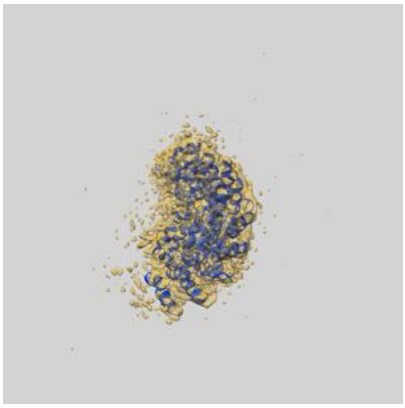
9.1 Map-model overlay



X



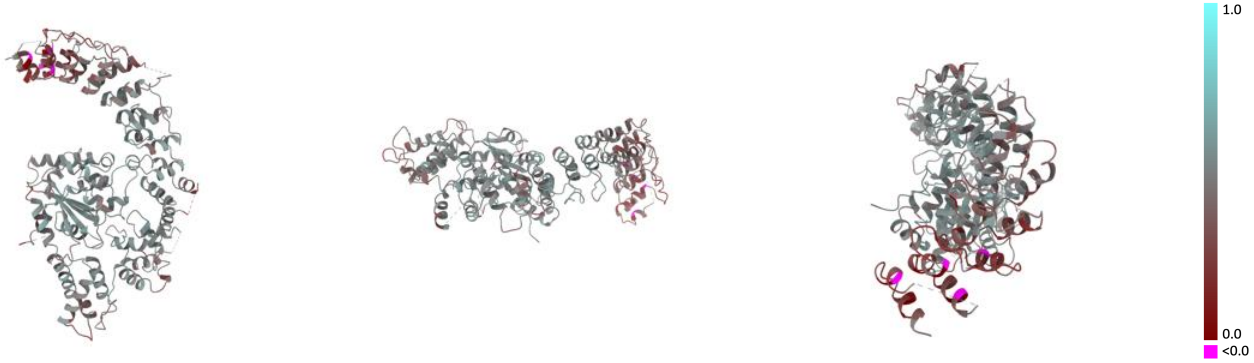
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0298 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.

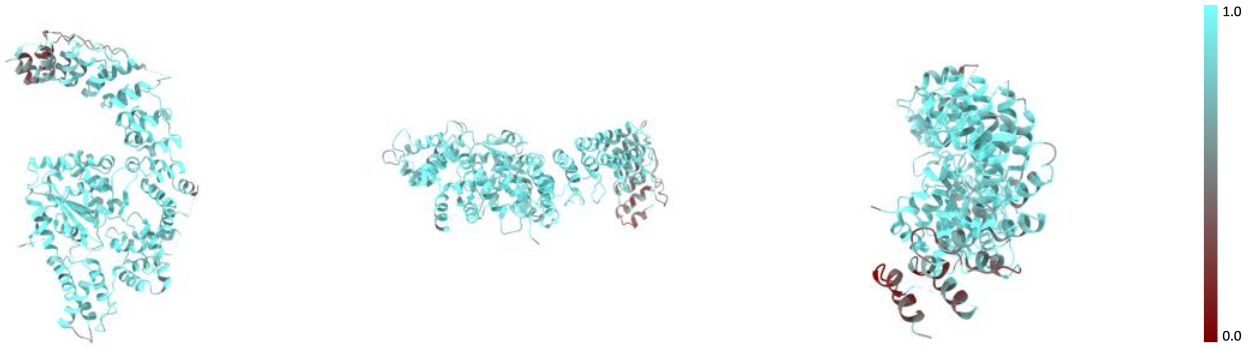
9.2 Q-score mapped to coordinate model (i)



X Y Z

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.

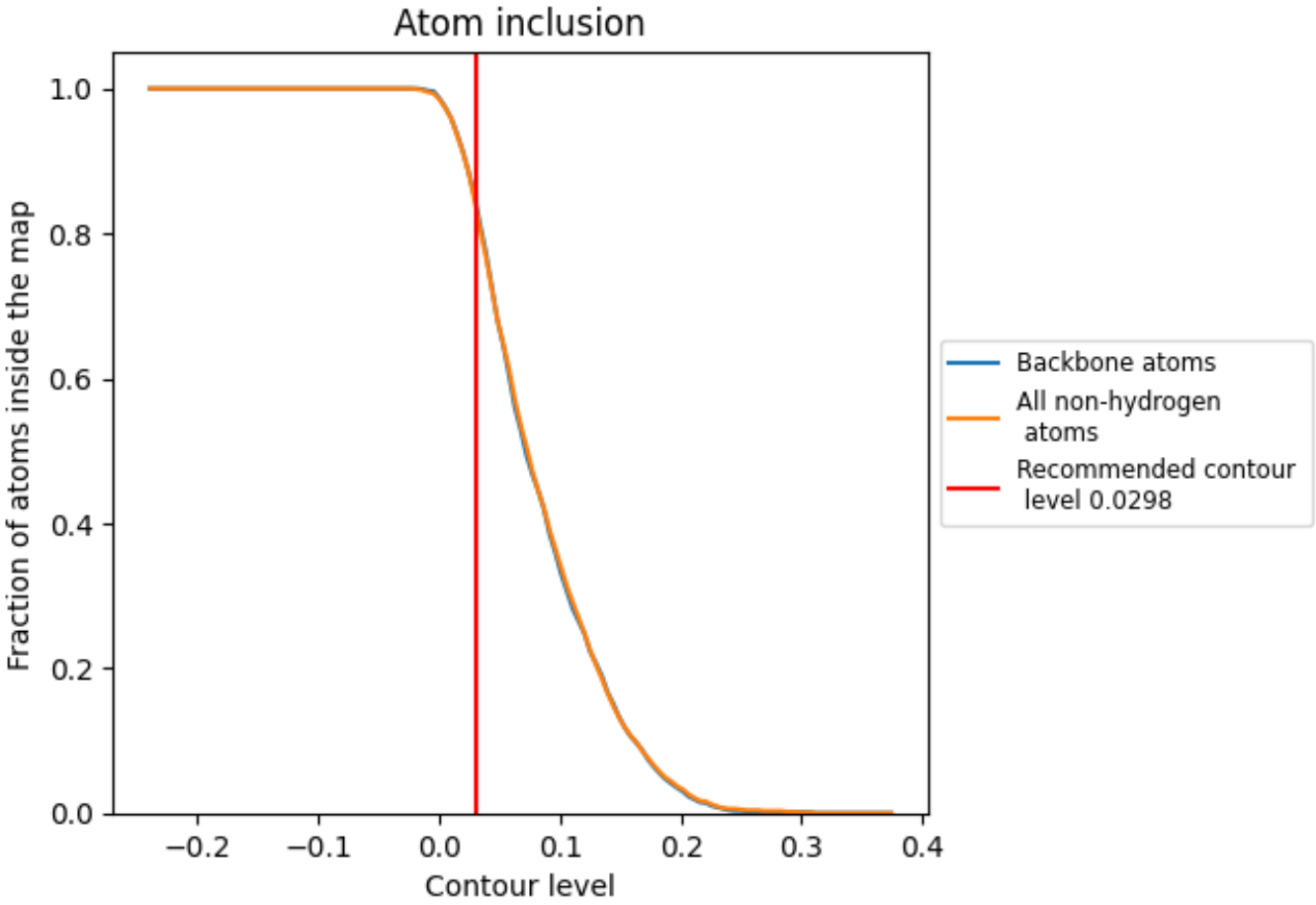
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.0298).

9.4 Atom inclusion (i)



At the recommended contour level, 85% of all backbone atoms, 84% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.0298) and Q-score for the entire model and for each chain.



Chain	Atom inclusion	Q-score
All	0.8420	0.4520
A	0.8550	0.4520