



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 04:08 AM UTC

PDB ID : 8FKY / pdb_00008fky
EMDB ID : EMD-29261
Title : Human nucleolar pre-60S ribosomal subunit (State F)
Authors : Vanden Broeck, A.; Klinge, S.
Deposited on : 2022-12-21
Resolution : 2.67 Å(reported)

This is a Full wwPDB EM Validation 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 : **FAILED**
MolProbity : **FAILED**
Buster-report : **FAILED**
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

PERCENTILES INFOmissingINFO

ENTRY-COMPOSITION INFOmissingINFO

SEQUENCE-PLOTS INFOmissingINFO

1 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	76907	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	64000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	9.161	Depositor
Minimum map value	0.000	Depositor
Average map value	0.051	Depositor
Map value standard deviation	0.192	Depositor
Recommended contour level	0.85	Depositor
Map size (Å)	514.56, 514.56, 514.56	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.072, 1.072, 1.072	Depositor

2 Model quality [i](#)

2.1 Standard geometry [i](#)

MolProbity failed to run properly - this section is therefore empty.

2.2 Too-close contacts [i](#)

MolProbity failed to run properly - this section is therefore empty.

2.3 Torsion angles [i](#)

2.3.1 Protein backbone [i](#)

MolProbity failed to run properly - this section is therefore empty.

2.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

2.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

2.4 Non-standard residues in protein, DNA, RNA chains [i](#)

Mogul failed to run properly - this section is therefore empty.

2.5 Carbohydrates [i](#)

Mogul failed to run properly - this section is therefore empty.

2.6 Ligand geometry [i](#)

Mogul failed to run properly - this section is therefore empty.

2.7 Other polymers [i](#)

Mogul failed to run properly - this section is therefore empty.

2.8 Polymer linkage issues

There are no chain breaks in this entry.

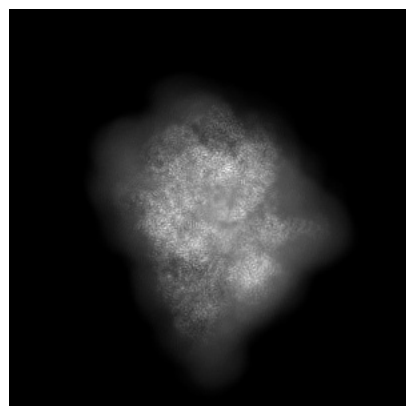
3 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29261. These allow visual inspection of the internal detail of the map and identification of artifacts.

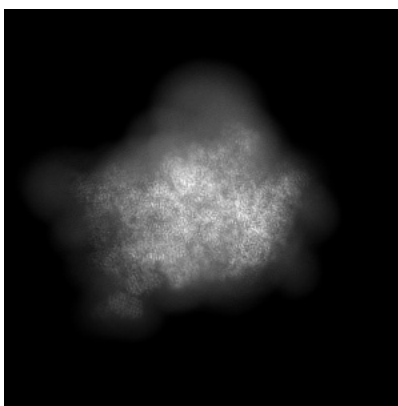
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

3.1 Orthogonal projections [i](#)

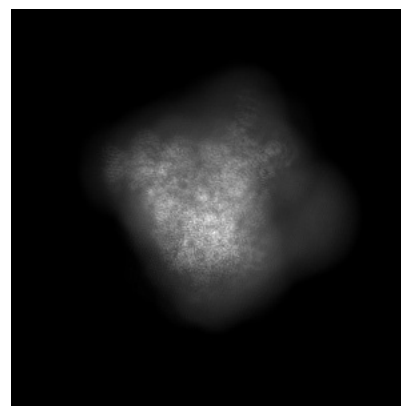
3.1.1 Primary map



X

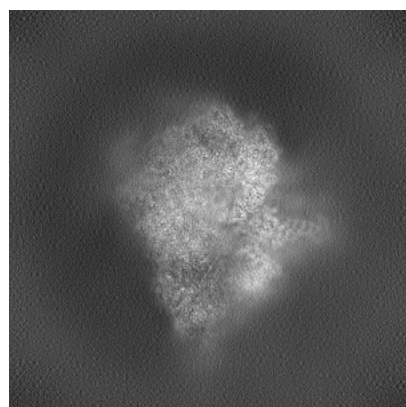


Y

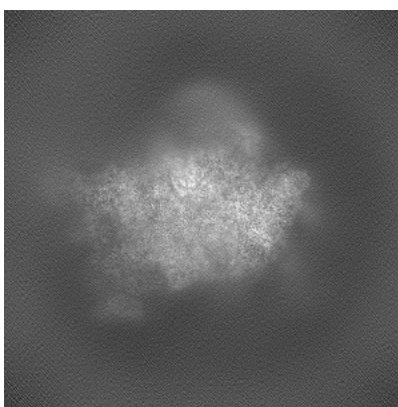


Z

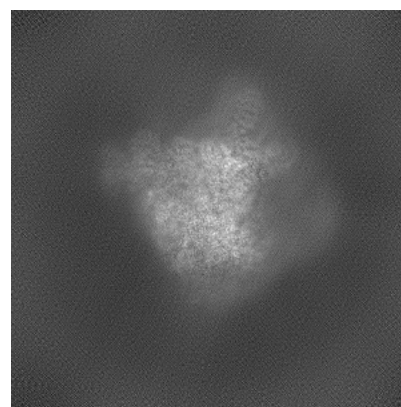
3.1.2 Raw map



X



Y

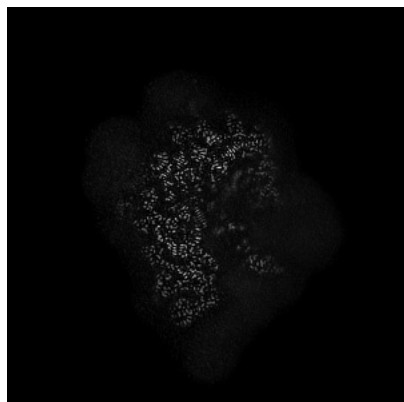


Z

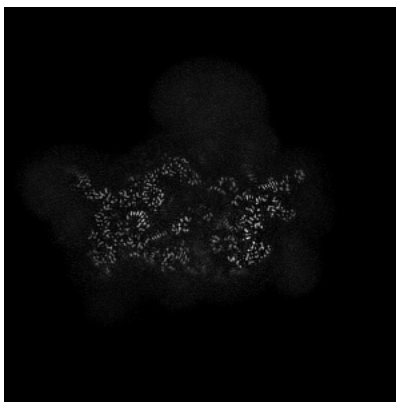
The images above show the map projected in three orthogonal directions.

3.2 Central slices [i](#)

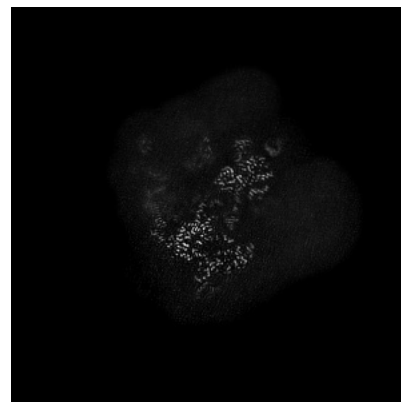
3.2.1 Primary map



X Index: 240

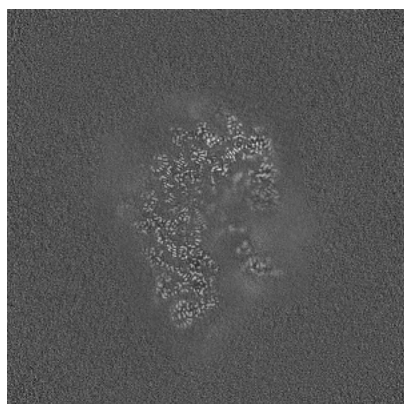


Y Index: 240

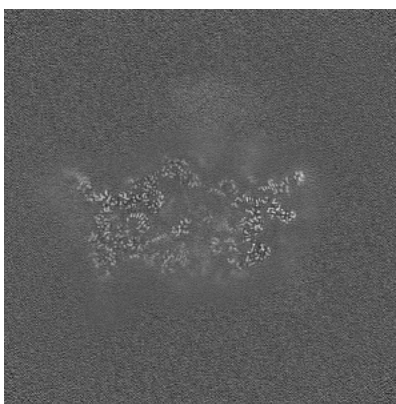


Z Index: 240

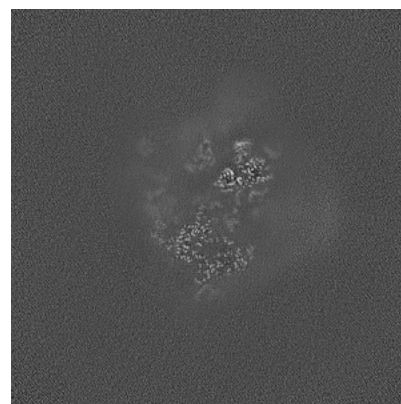
3.2.2 Raw map



X Index: 240



Y Index: 240

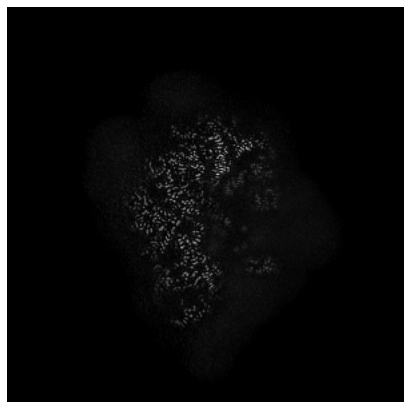


Z Index: 240

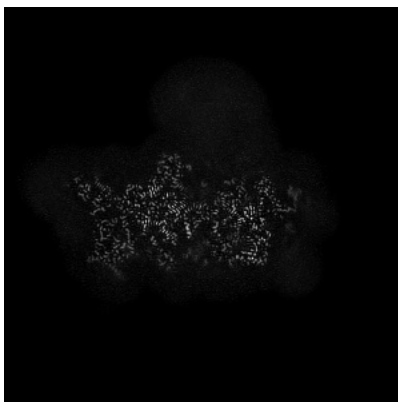
The images above show central slices of the map in three orthogonal directions.

3.3 Largest variance slices [i](#)

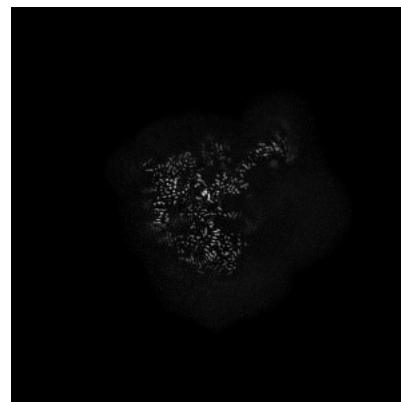
3.3.1 Primary map



X Index: 233

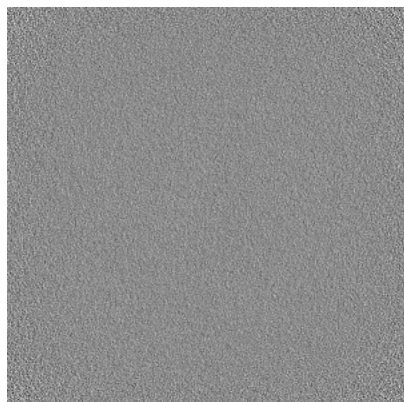


Y Index: 228

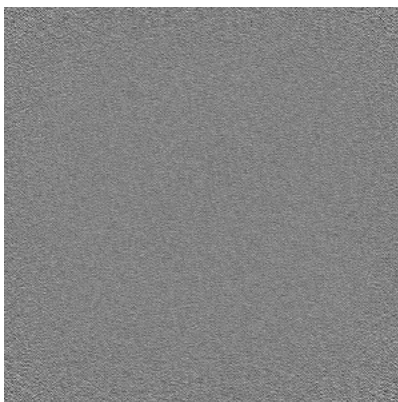


Z Index: 285

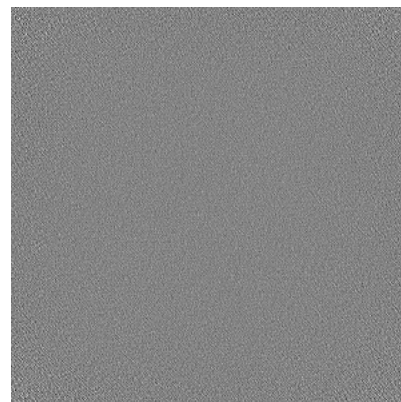
3.3.2 Raw map



X Index: 0



Y Index: 0

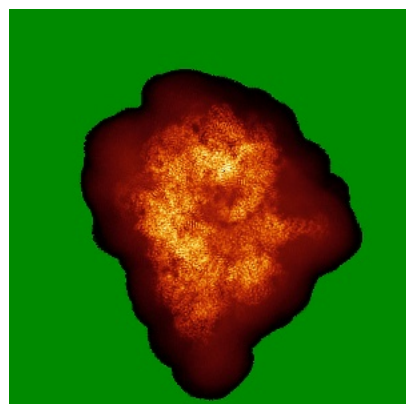


Z Index: 0

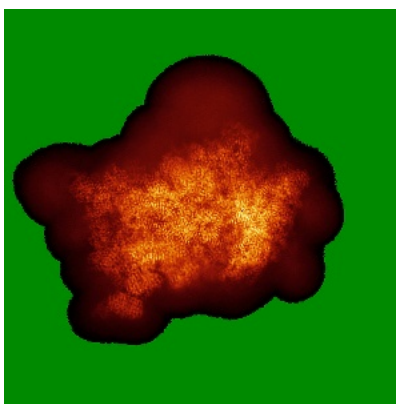
The images above show the largest variance slices of the map in three orthogonal directions.

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

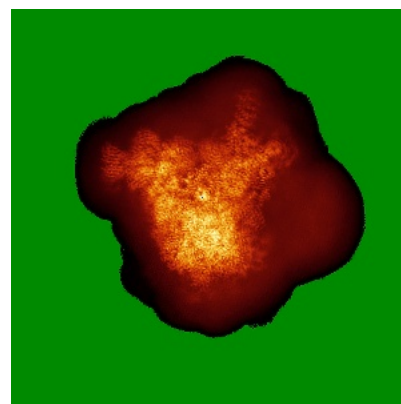
3.4.1 Primary map



X

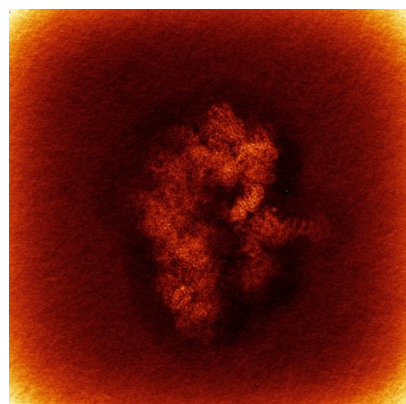


Y

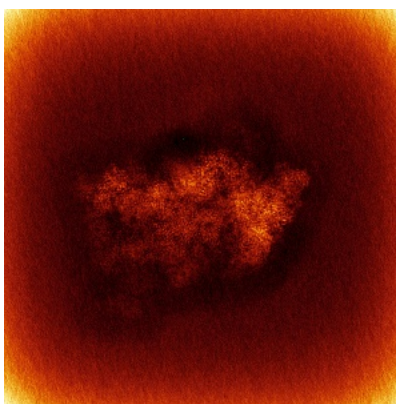


Z

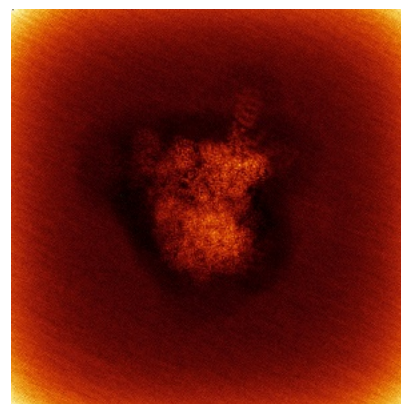
3.4.2 Raw map



X



Y



Z

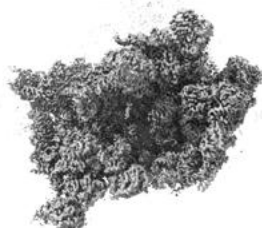
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

3.5 Orthogonal surface views [i](#)

3.5.1 Primary map



X



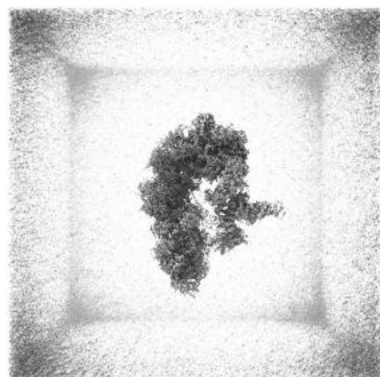
Y



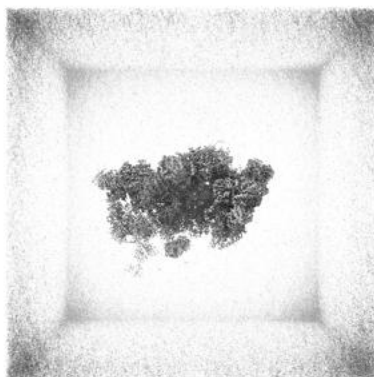
Z

The images above show the 3D surface view of the map at the recommended contour level 0.85. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

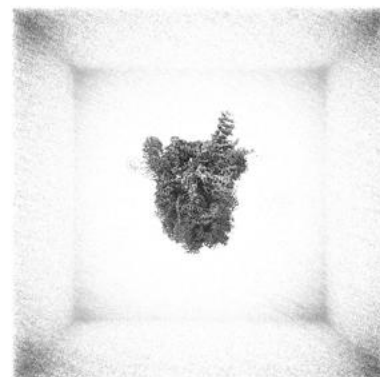
3.5.2 Raw map



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.

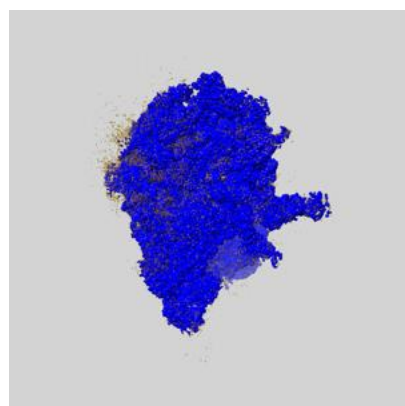
3.6 Mask visualisation [i](#)

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

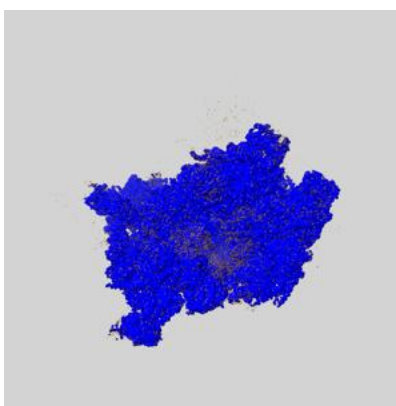
A mask typically either:

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

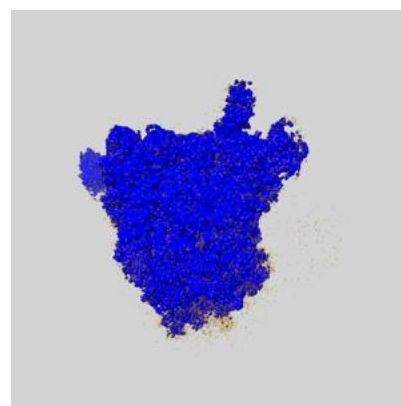
3.6.1 emd_29261_msk_1.map [i](#)



X



Y

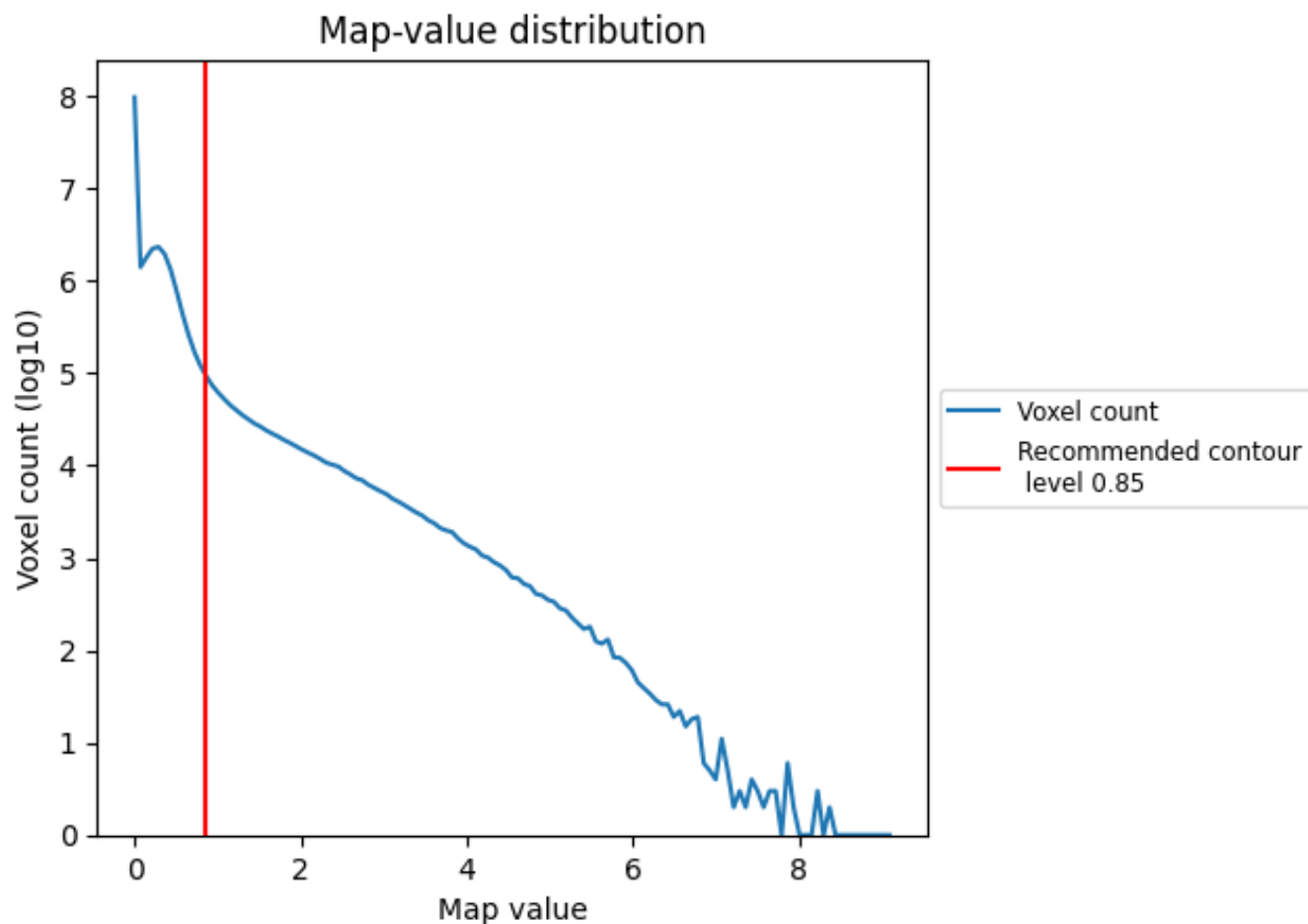


Z

4 Map analysis [i](#)

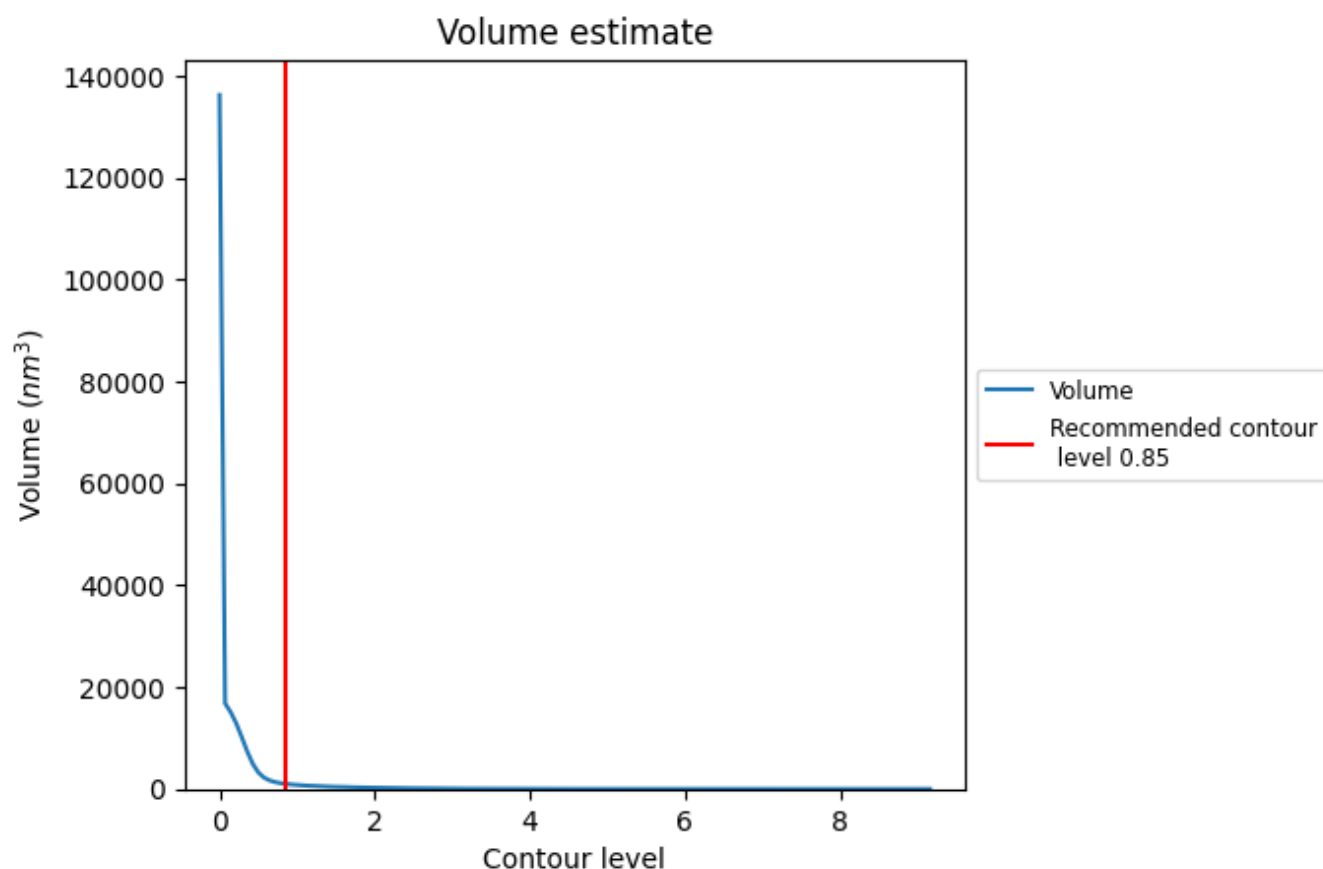
This section contains the results of statistical analysis of the map.

4.1 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.

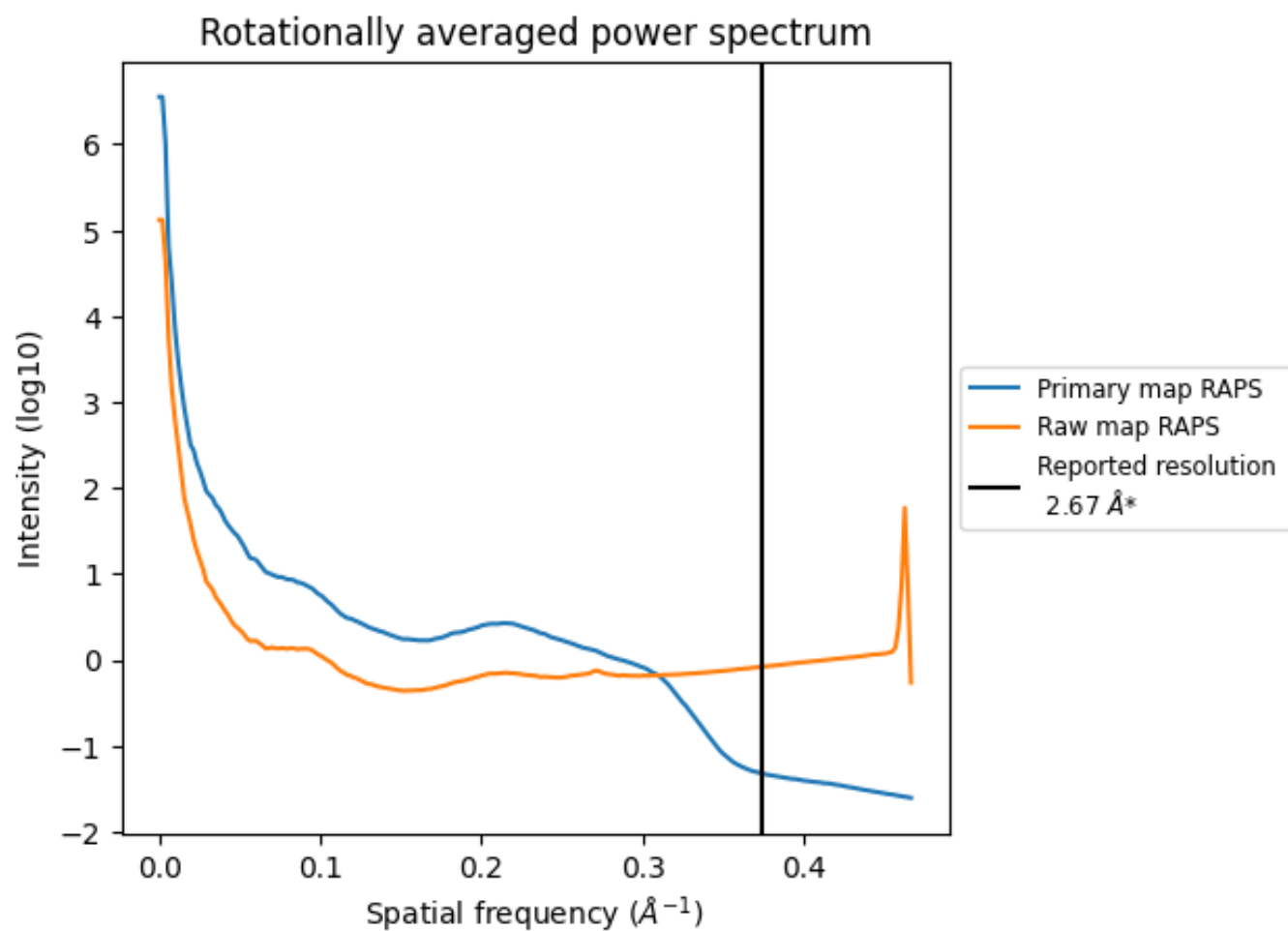
4.2 Volume estimate [i](#)



The volume at the recommended contour level is 1004 nm^3 ; this corresponds to an approximate mass of 907 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.

4.3 Rotationally averaged power spectrum ⓘ

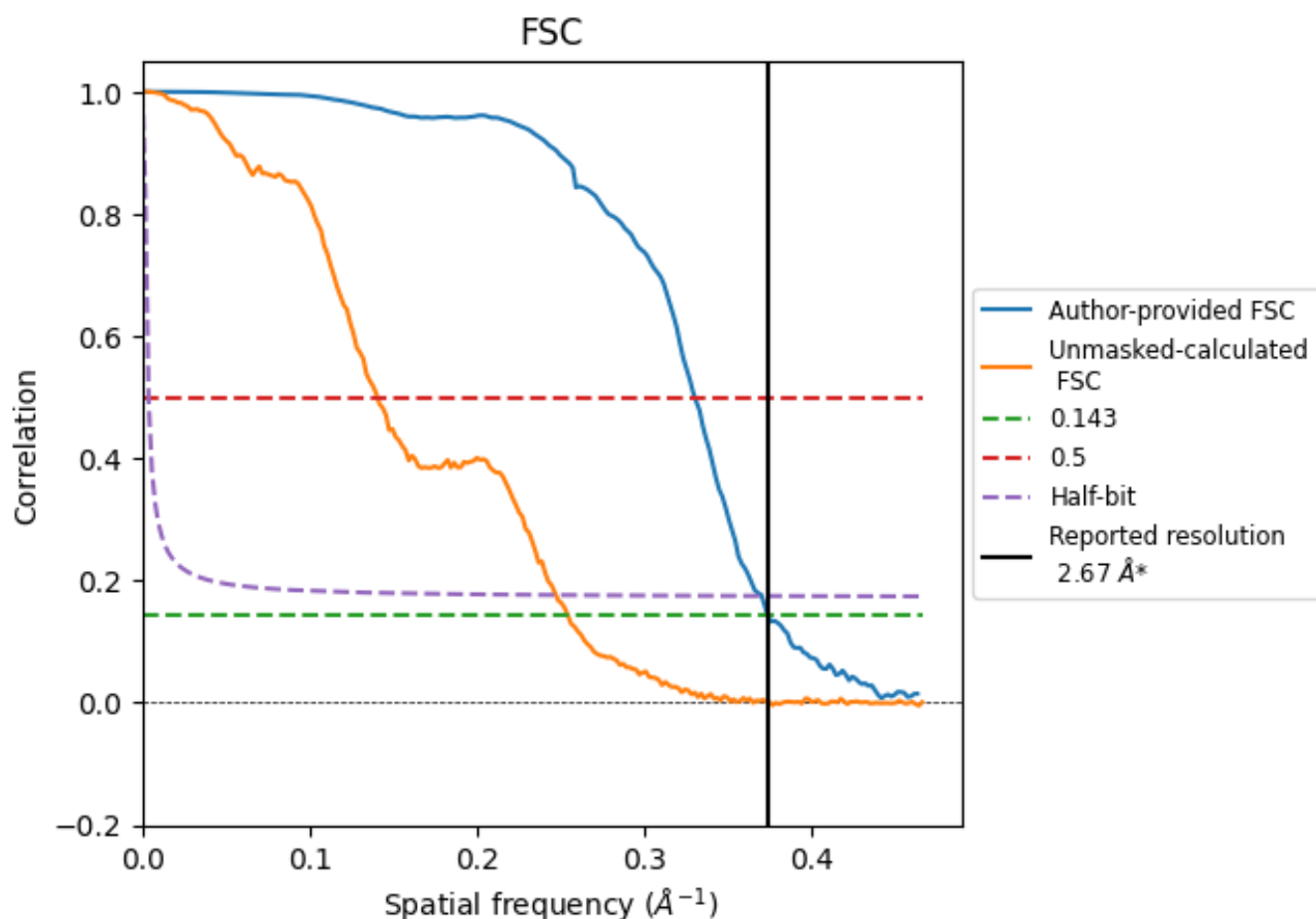


*Reported resolution corresponds to spatial frequency of 0.375 $\text{\AA}^{-1}$

5 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.

5.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.375 $\text{\AA}^{-1}$

5.2 Resolution estimates [i](#)

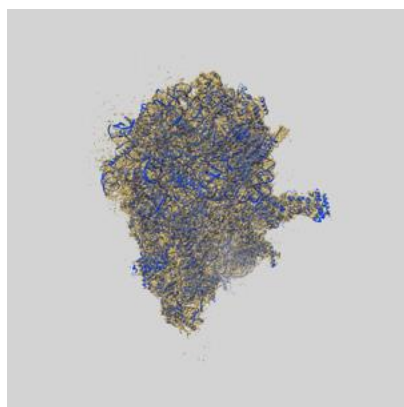
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.67	-	-
Author-provided FSC curve	2.67	3.03	2.70
Unmasked-calculated*	3.93	7.11	4.04

*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 3.93 differs from the reported value 2.67 by more than 10 %

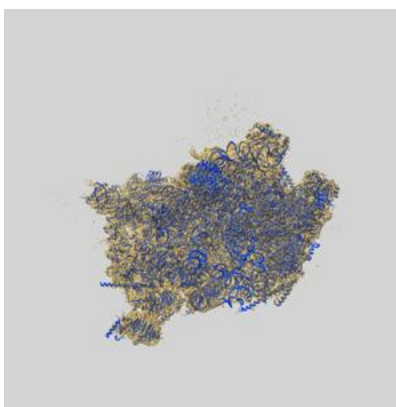
6 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29261 and PDB model 8FKY. Per-residue inclusion information can be found in section ?? on page ??.

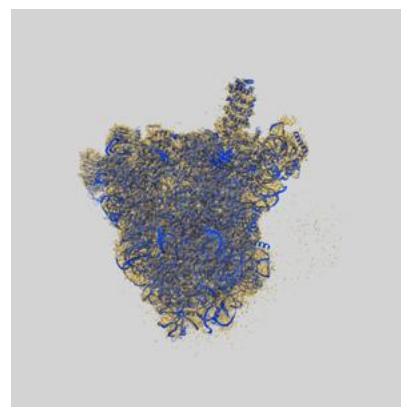
6.1 Map-model overlay [i](#)



X



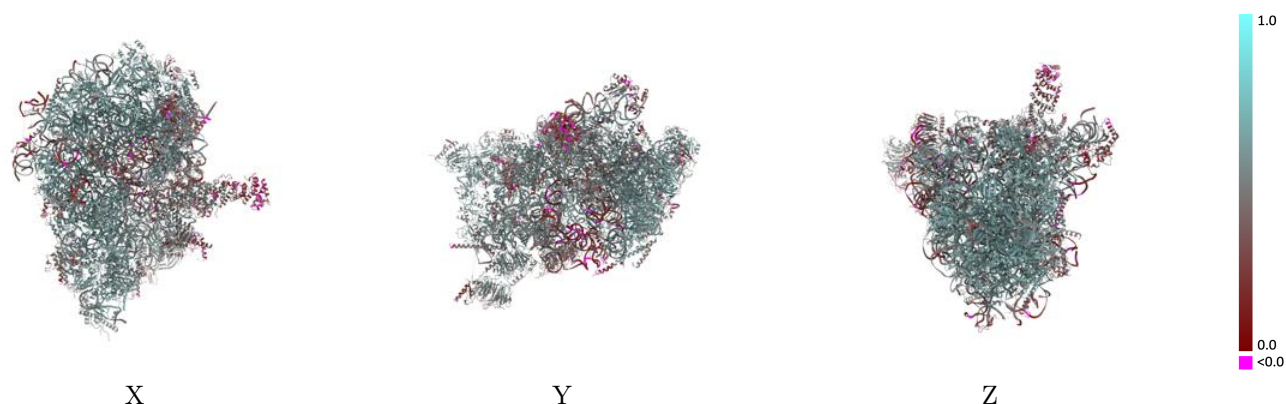
Y



Z

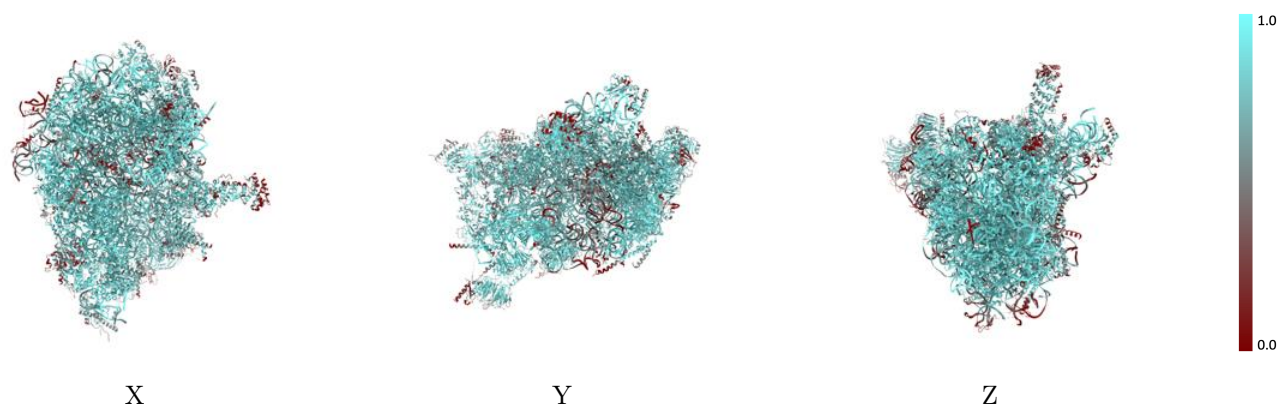
The images above show the 3D surface view of the map at the recommended contour level 0.85 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



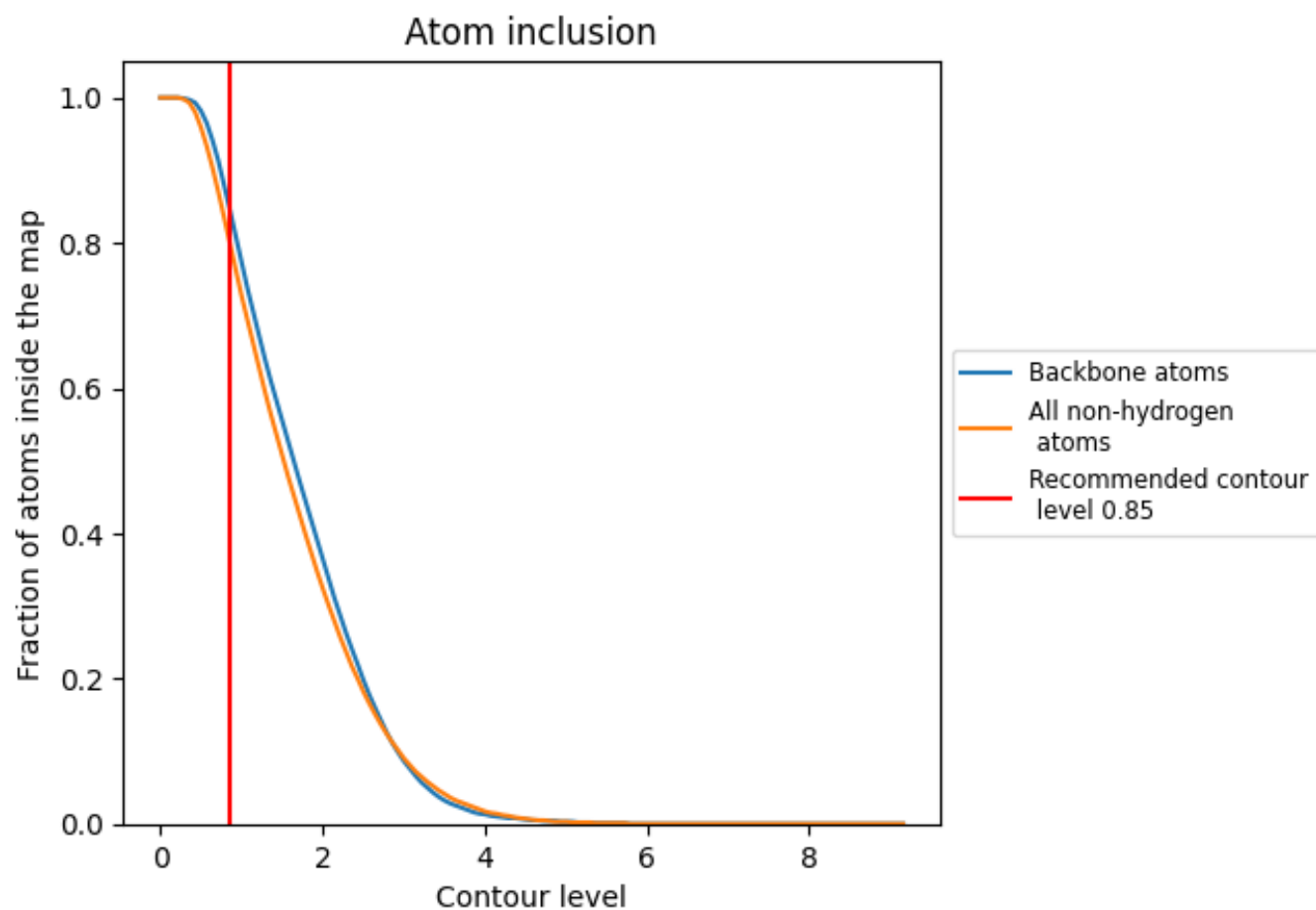
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

6.3 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 (0.85).

6.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 80% of all non-hydrogen atoms, are inside the map.

6.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.85) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8040	 0.5190
BA	 0.4420	 0.4340
BB	 0.6010	 0.3170
BC	 0.7330	 0.4050
L1	 0.9340	 0.5760
L2	 0.8130	 0.5490
L3	 0.8430	 0.5110
L6	 0.9330	 0.6060
L7	 0.9020	 0.6130
L8	 0.8620	 0.6050
L9	 0.9690	 0.6290
LA	 0.8080	 0.5590
LB	 0.9380	 0.6140
LC	 0.8790	 0.6200
LD	 0.6660	 0.3870
LE	 0.4810	 0.4200
LG	 0.8330	 0.5550
LH	 0.8110	 0.5360
LI	 0.8790	 0.6020
LJ	 0.8820	 0.5040
LK	 0.6990	 0.4340
LL	 0.9120	 0.6030
LN	 0.9090	 0.6100
LO	 0.8410	 0.4810
LP	 0.2800	 0.2770
LQ	 0.9420	 0.6280
LR	 0.8380	 0.4760
LS	 0.8820	 0.5770
LT	 0.9510	 0.6310
LU	 0.8310	 0.5830
LW	 0.9490	 0.6110
NA	 0.7250	 0.3810
NB	 0.6680	 0.5320
NF	 0.8440	 0.5740
NH	 0.8950	 0.5450



Continued on next page...

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Chain	Atom inclusion	Q-score
NI	 0.5420	 0.4370
NK	 0.7860	 0.5560
SA	 0.9220	 0.6140
SC	 0.5950	 0.4930
SD	 0.8660	 0.5870
SE	 0.9060	 0.5940
SG	 0.8860	 0.6090
SH	 0.7130	 0.5480
SI	 0.7840	 0.5610
SJ	 0.6090	 0.4250
SK	 0.8560	 0.5630
SL	 0.7300	 0.5450
SM	 0.8170	 0.5310
SN	 0.7080	 0.4910
SO	 0.7850	 0.5440
SP	 0.6430	 0.4300
SQ	 0.7750	 0.5660
SR	 0.8090	 0.5520
SS	 0.8490	 0.5220
ST	 0.5830	 0.4000
SU	 0.7210	 0.4380
SV	 0.7510	 0.5140
SW	 0.6480	 0.5030
SY	 0.8740	 0.5460
SZ	 0.8460	 0.5580