



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 12:13 PM UTC

PDB ID : 7UIH / pdb_00007uih
EMDB ID : EMD-24743
Title : PSMD2 Structure
Authors : Johnson, M.C.; Bashore, C.; Ciferri, C.; Dueber, E.C.
Deposited on : 2022-03-29
Resolution : 3.10 Å(reported)
Based on initial model : 6MSJ

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:

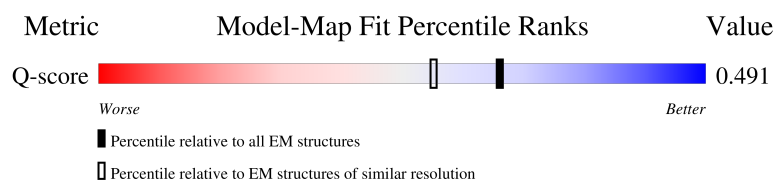
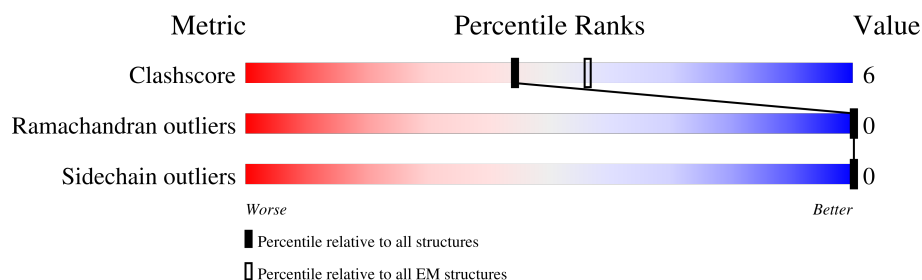
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.10 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	14724 (2.60 - 3.60)

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	A	644	11% 75% 15% 11%
2	C	217	9% 90%
3	D	231	11% 89%
4	E	217	7% 92%

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Mol	Chain	Length	Quality of chain
5	F	229	8% • 90%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5114 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 26S proteasome non-ATPase regulatory subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	574	Total	C	N	O	S	0	0
			4379	2782	741	826	30		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	469	PHE	TYR	conflict	UNP Q13200

- Molecule 2 is a protein called Fab 14 LC CDRs.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	C	21	Total	C	N	O	0	0
			172	111	24	37		

- Molecule 3 is a protein called Fab 14 HC CDRs.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	D	26	Total	C	N	O	0	0
			223	147	32	44		

- Molecule 4 is a protein called Fab 8 LC CDRs.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	E	18	Total	C	N	O	0	0
			134	85	18	31		

- Molecule 5 is a protein called Fab 8 HC CDRs.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	23	Total	C	N	O	S	0	0
			206	137	31	37	1		

LYS
HIS
LYS
VAL
TYR
CYS
GLU
GLY
HIS
THR
GLY
GLY
LEU
SER
SER
VAL
THR
LYS
SER
PHE
ASN
ARG
GLY
GLU
CYS

● Molecule 3: Fab 14 HC CDRs

Chain D: 11% 89%

GLU
ILE
SER
GLU
VAL
GLN
LEU
VAL
GLU
SER
GLY
HIS
GLY
LEU
VAL
GLN
PRO
MET
GLY
GLY
SER
LEU
VAL
GLN
SER
PHE
ARG
LEU
SER
CYS
ALA
VAL
TYR
Y34
H38
TRP
VAL
ARG
GLN
ALA
PRO
GLY
LYS
GLY
LEU
TRP
VAL
452
Y58
S62
TYR
ALA
ASP
SER
VAL
LYS

GLY
ARG
PHE
THR
ILE
SER
ALA
ASP
THR
GLY
SER
LYS
ASN
GLY
ALA
TYR
GLN
MET
ASN
SER
LEU
ARG
ALA
GLU
SER
CYS
ALA
THR
THR
VAL
VAL
TYR
TRP
TYR
CYS
ASN
GLY
R101
G110
LEU
ASP
TYR
TRP
GLY
GLN
THR
GLY
LEU
VAL
VAL
SER
TRP
LYS
ASN
SER
GLY
ASP
LYS
ALA
LEU
THR
THR
HIS

ALA
PRO
SER
SER
SER
LEU
GLY
THR
SER
GLY
THR
GLY
THR
ILE
CYS
ALA
LEU
GLY
CYS
HIS
LYS
LEU
VAL
PRO
VAL
LYS
ASP
SER
ASN
THR
PHE
PRO
VAL
PRO
GLY
VAL
THR
SER
GLU
PRO
LYS
CYS
SER
ASN
GLY
ASP
LYS
THR
HIS

● Molecule 4: Fab 8 LC CDRs

Chain E: 7% 92%

SER
ASP
ILE
SER
MET
THR
GLN
SER
PRO
SER
SER
LEU
SER
ALA
SER
VAL
GLN
ASP
ARG
VAL
THR
ILE
THR
THR
CYS
ALA
SER
GLN
GLN
Y30
Y34
ALA
TRP
TYR
GLN
GLN
LYS
PRO
GLY
LYS
VAL
ALA
PRO
LYS
LEU
ILE
LEU
ILE
Y50
S54
LEU
TYR
SER
SER
GLY
VAL
PRO
SER
ARG
PHE
SER

GLY
SER
ARG
SER
GLY
THR
ASP
PHE
THR
LEU
SER
ILE
SER
LEU
ASN
VAL
GLN
PRO
GLU
ASP
PHE
ALA
THR
ILE
THR
TYR
CYS
GLN
GLN
Y92
S96
S97
L98
VAL
THR
PHE
GLY
ASN
GLY
GLN
GLY
SER
GLY
VAL
THR
GLU
ILE
LYS
ASP
THR
GLY
GLY
CYS
GLU
LYS

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

ALA
ASP
TYR
GLY
HIS
LYS
VAL
TYR
ALA
CYS
GLY
VAL
THR
HIS
GLN
GLY
SER
SER
GLY
PRO
VAL
THR
LYS
SER
PHE
ASN
GLY
GLY
CYS

● Molecule 5: Fab 8 HC CDRs

Chain F: 8% 90%

GLU
ILE
SER
GLU
VAL
GLN
LEU
VAL
GLU
SER
GLY
HIS
GLY
LEU
VAL
GLN
PRO
GLY
SER
LEU
VAL
GLN
MET
ARG
SER
CYS
ALA
ALA
SER
GLY
PHE
ASN
PHE
S33
H38
TRP
VAL
ARG
GLN
ALA
LEU
PRO
GLY
LYS
GLY
LEU
TRP
VAL
Y53
I54
Y55
P56
Y57
S58
S59
Y62
ALA

ASP
SER
VAL
PHE
PRO
GLY
ARG
PHE
THR
ILE
SER
ALA
ASP
THR
SER
GLY
VAL
ASN
THR
ALA
GLY
LEU
TYR
LEU
GLN
MET
SER
SER
ARG
ALA
SER
GLY
ASP
THR
ALA
VAL
VAL
TYR
TYR
CYS
ALA
ARG
K102
A108
LEU
ASP
SER
TYR
GLY
GLY
GLN
THR
THR
VAL
VAL
THR
VAL
GLN
SER
SER
SER
GLY
ALA
SER
THR
LYS
GLY
PRO

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

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	232101	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 ($e^-/\text{\AA}^2$)	64	Depositor
Minimum defocus (nm)	300	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.025	Depositor
Minimum map value	-0.695	Depositor
Average map value	-0.002	Depositor
Map value standard deviation	0.039	Depositor
Recommended contour level	0.2	Depositor
Map size ($\text{\AA}$)	402.24, 402.24, 402.24	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.838, 0.838, 0.838	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.13	0/4448	0.39	0/6026
2	C	0.15	0/175	0.28	0/236
3	D	0.12	0/231	0.33	0/311
4	E	0.11	0/135	0.38	0/180
5	F	0.17	0/214	0.39	0/287
All	All	0.14	0/5203	0.39	0/7040

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4379	0	4435	61	0
2	C	172	0	153	1	0
3	D	223	0	188	1	0
4	E	134	0	117	2	0
5	F	206	0	178	4	0
All	All	5114	0	5071	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 6.

All (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:A:392:THR:HG22	1:A:393:ASP:H	1.41	0.85
1:A:283:THR:HG22	1:A:284:SER:H	1.44	0.81
1:A:710:LEU:O	1:A:714:SER:OG	2.00	0.79
4:E:96:SER:O	4:E:97:SER:OG	2.03	0.75
1:A:680:ARG:O	2:C:94:SER:OG	2.03	0.75
1:A:829:MET:HE1	1:A:831:VAL:CG2	2.26	0.65
1:A:581:GLU:O	1:A:588:ARG:NH2	2.30	0.65
5:F:58:SER:O	5:F:59:SER:OG	2.15	0.65
1:A:333:LEU:HD12	1:A:871:PRO:HB2	1.80	0.64
1:A:608:LYS:O	1:A:612:LEU:HD23	1.99	0.63
1:A:388:ASP:OD2	1:A:391:LEU:N	2.29	0.63
1:A:341:GLU:O	1:A:773:LYS:NZ	2.35	0.60
1:A:445:LEU:HD22	1:A:466:LEU:HD23	1.85	0.59
1:A:834:ASP:OD1	1:A:835:GLU:N	2.35	0.57
1:A:392:THR:HG22	1:A:393:ASP:N	2.16	0.57
1:A:741:LEU:O	1:A:745:LEU:HD23	2.06	0.54
1:A:534:VAL:HG13	1:A:565:ASN:ND2	2.23	0.53
1:A:829:MET:HE1	1:A:831:VAL:HG21	1.90	0.52
1:A:846:VAL:HG23	1:A:846:VAL:O	2.10	0.52
1:A:870:THR:HB	1:A:871:PRO:HA	1.92	0.52
1:A:372:LEU:HD11	1:A:376:PHE:CE2	2.44	0.51
1:A:283:THR:HG22	1:A:284:SER:N	2.21	0.51
1:A:505:MET:HE1	1:A:538:ILE:HD13	1.92	0.51
1:A:740:ARG:HH11	1:A:740:ARG:HG2	1.76	0.50
1:A:431:LYS:HB3	5:F:55:TYR:OH	2.12	0.49
4:E:30:VAL:HG22	4:E:30:VAL:O	2.11	0.49
5:F:55:TYR:CE2	5:F:57:TYR:HB3	2.48	0.48
1:A:659:LEU:HD12	1:A:662:MET:HE2	1.96	0.48
1:A:547:GLU:OE1	1:A:547:GLU:N	2.45	0.48
1:A:478:ARG:NH1	1:A:507:ASP:OD2	2.41	0.47
1:A:829:MET:CE	1:A:831:VAL:CG2	2.93	0.47
1:A:545:LYS:HB2	1:A:550:LEU:HD21	1.97	0.47
1:A:416:MET:HE3	1:A:822:VAL:HG12	1.97	0.46
1:A:653:ALA:O	1:A:657:ILE:HG13	2.15	0.46
1:A:333:LEU:C	1:A:333:LEU:HD23	2.42	0.45
1:A:531:ASN:HB3	1:A:534:VAL:HG12	1.98	0.45
1:A:683:GLU:OE2	3:D:58:TYR:CD1	2.69	0.45
1:A:367:SER:N	1:A:370:MET:HE2	2.31	0.45
1:A:590:PHE:CE1	1:A:652:VAL:HG21	2.51	0.45
1:A:844:VAL:HA	1:A:882:LEU:HA	1.99	0.45
1:A:431:LYS:CB	5:F:55:TYR:OH	2.65	0.45
1:A:373:ALA:O	1:A:377:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:ASP:O	1:A:394:ASP:OD1	2.35	0.45
1:A:441:LYS:HG2	1:A:477:MET:HE1	1.99	0.45
1:A:760:PHE:CE2	1:A:764:LEU:HD11	2.52	0.45
1:A:716:ASP:OD1	1:A:717:ALA:N	2.50	0.44
1:A:749:ALA:HB2	1:A:762:VAL:HG11	2.00	0.43
1:A:895:GLU:HA	1:A:895:GLU:OE1	2.19	0.43
1:A:305:LEU:O	1:A:305:LEU:HD12	2.19	0.42
1:A:563:GLY:O	1:A:567:LEU:N	2.52	0.42
1:A:266:LEU:O	1:A:270:LEU:HG	2.20	0.41
1:A:872:VAL:HG22	1:A:873:LEU:N	2.35	0.41
1:A:606:VAL:O	1:A:610:GLN:HG2	2.20	0.41
1:A:458:GLU:O	1:A:458:GLU:HG3	2.20	0.41
1:A:372:LEU:HD11	1:A:376:PHE:HE2	1.84	0.41
1:A:573:ILE:HG23	1:A:577:LEU:HD13	2.03	0.41
1:A:408:LEU:C	1:A:408:LEU:HD23	2.45	0.41
1:A:322:SER:O	1:A:455:VAL:O	2.39	0.41
1:A:422:VAL:HG11	1:A:457:ASN:ND2	2.36	0.41
1:A:655:LEU:HD12	1:A:804:LEU:HD11	2.03	0.41
1:A:690:VAL:N	1:A:691:PRO:HD2	2.36	0.40
1:A:883:ALA:O	1:A:884:THR:CG2	2.69	0.40
1:A:671:ALA:O	1:A:675:PHE:CD2	2.75	0.40
1:A:819:TYR:O	1:A:822:VAL:HG22	2.21	0.40
1:A:474:SER:OG	1:A:476:THR:HG22	2.22	0.40

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	566/644 (88%)	554 (98%)	12 (2%)	0	100	100
2	C	15/217 (7%)	13 (87%)	2 (13%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	D	20/231 (9%)	19 (95%)	1 (5%)	0	100	100
4	E	12/217 (6%)	8 (67%)	4 (33%)	0	100	100
5	F	17/229 (7%)	15 (88%)	2 (12%)	0	100	100
All	All	630/1538 (41%)	609 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	476/534 (89%)	476 (100%)	0	100	100
2	C	20/192 (10%)	20 (100%)	0	100	100
3	D	23/194 (12%)	23 (100%)	0	100	100
4	E	16/192 (8%)	16 (100%)	0	100	100
5	F	21/192 (11%)	21 (100%)	0	100	100
All	All	556/1304 (43%)	556 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	327	ASN
1	A	565	ASN
1	A	752	HIS
1	A	848	GLN
5	F	38	HIS

5.3.3 RNA ⓘ

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](#)

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.

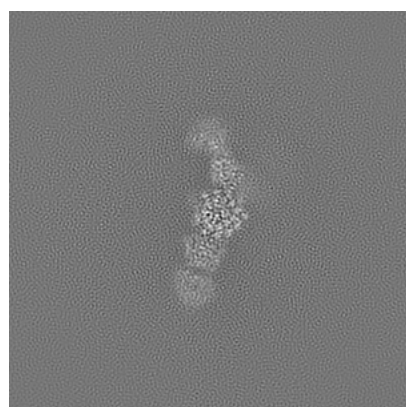
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24743. These allow visual inspection of the internal detail of the map and identification of artifacts.

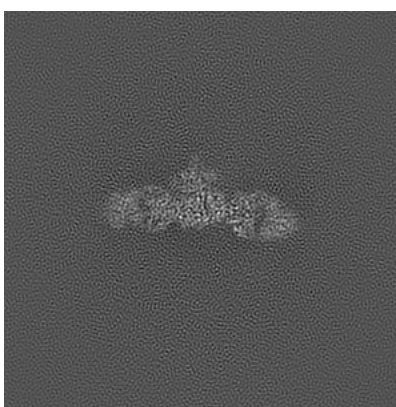
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

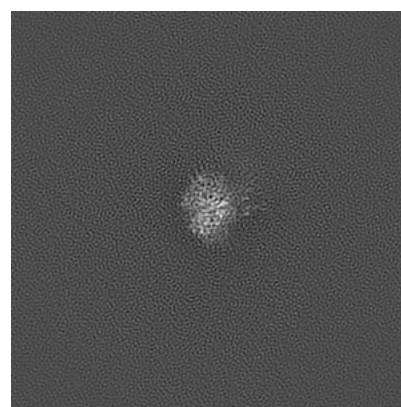
6.1.1 Primary map



X



Y

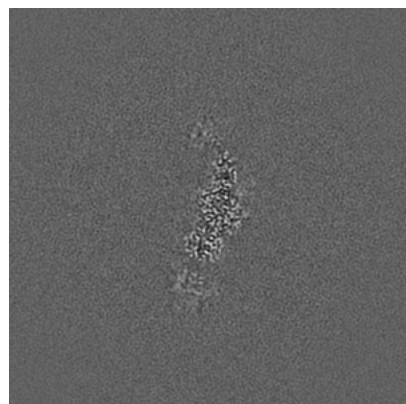


Z

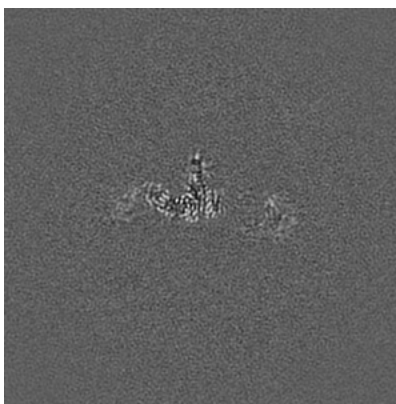
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

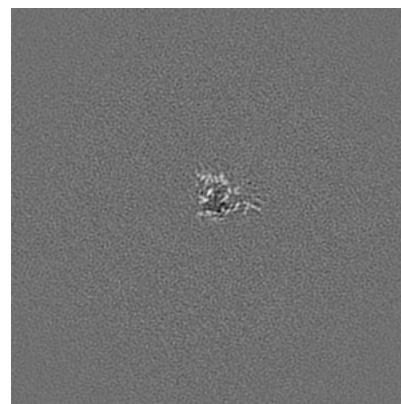
6.2.1 Primary map



X Index: 240



Y Index: 240

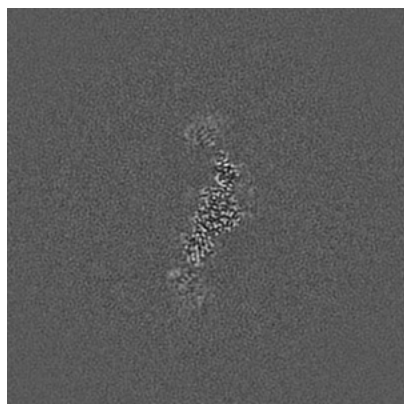


Z Index: 240

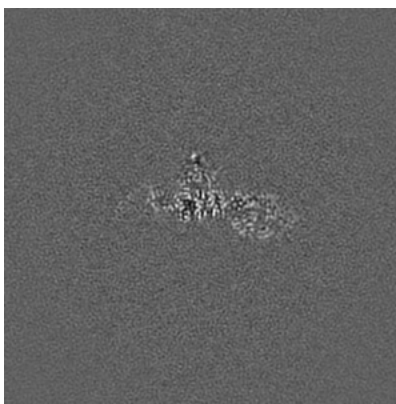
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

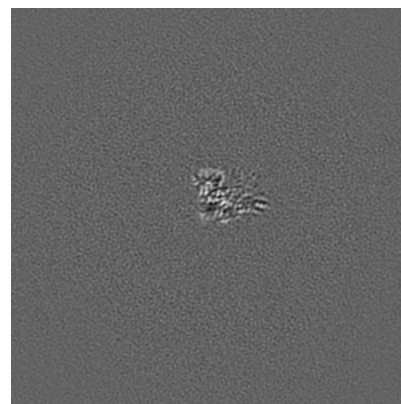
6.3.1 Primary map



X Index: 235



Y Index: 248

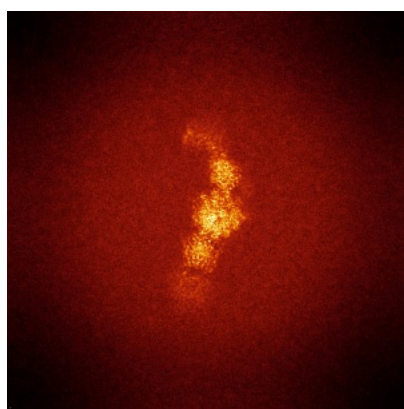


Z Index: 233

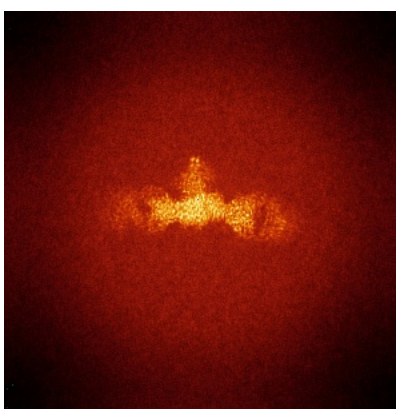
The images above show the largest variance slices of the map in three orthogonal directions.

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

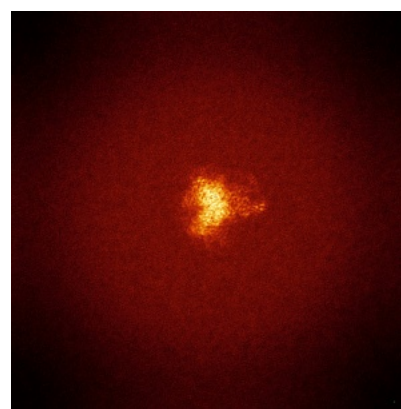
6.4.1 Primary map



X



Y



Z

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

This section was not generated.

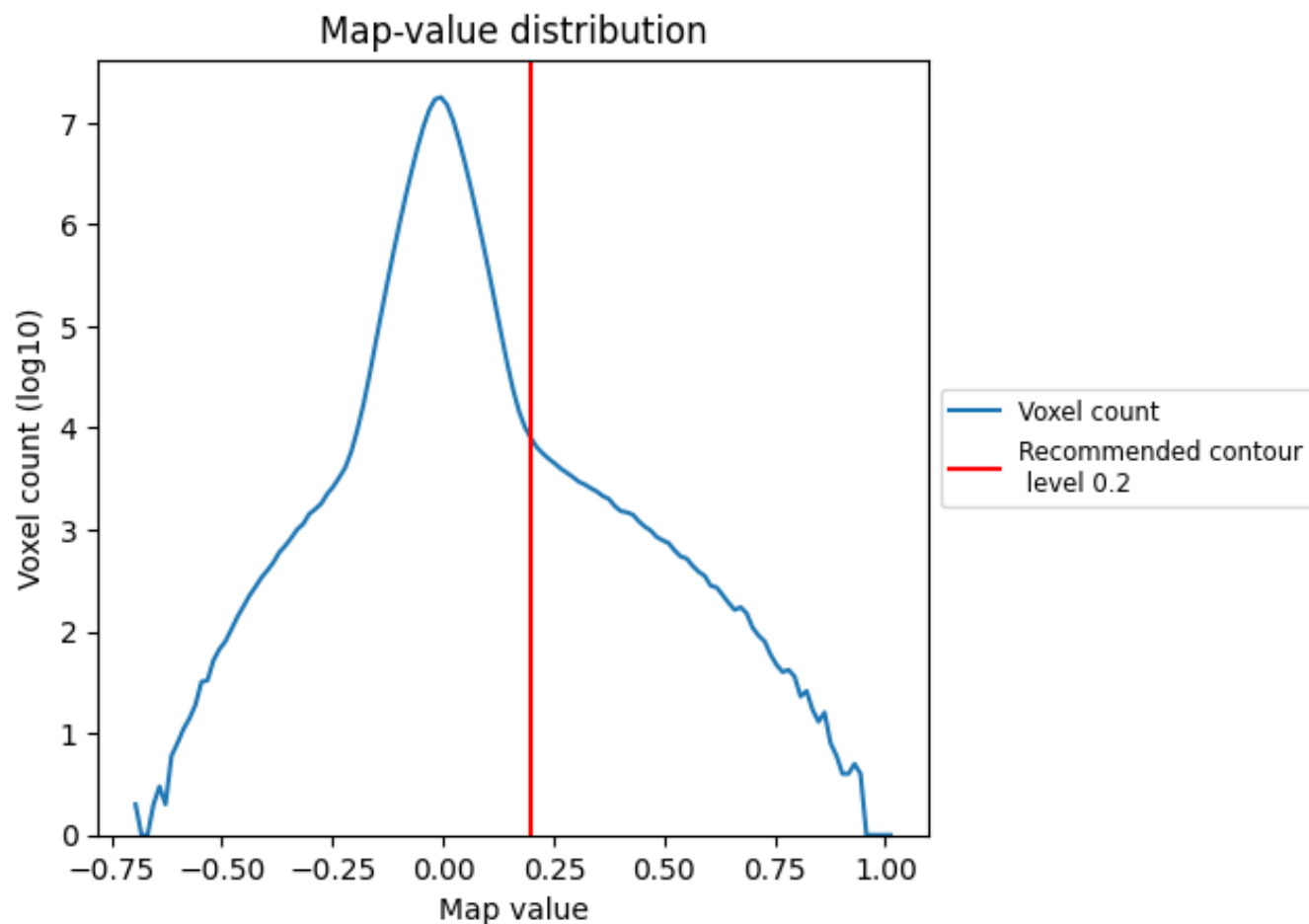
6.6 Mask visualisation

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

7 Map analysis [i](#)

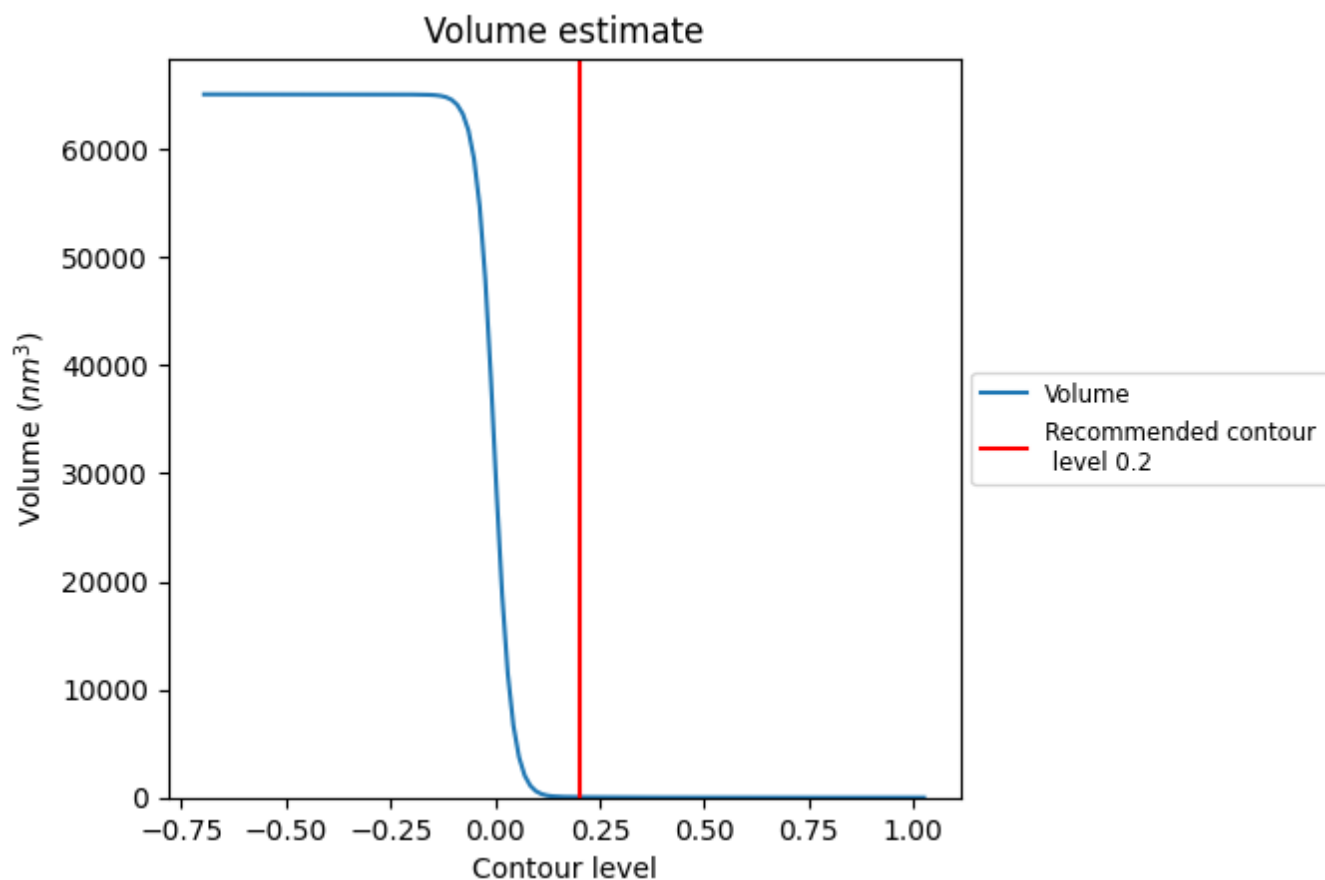
This section contains the results of statistical analysis of the map.

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.

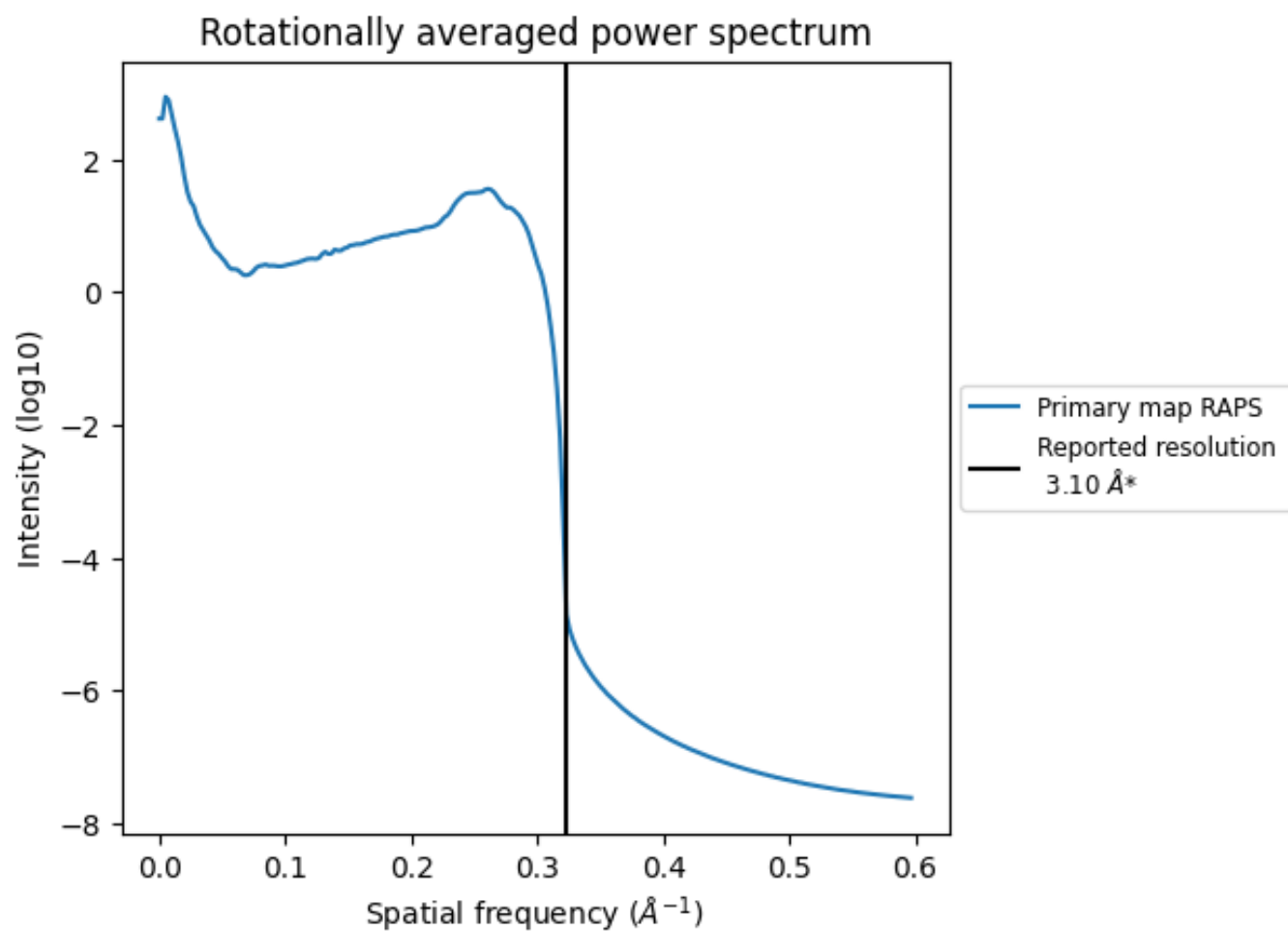
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 42 nm³; this corresponds to an approximate mass of 38 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 ⓘ

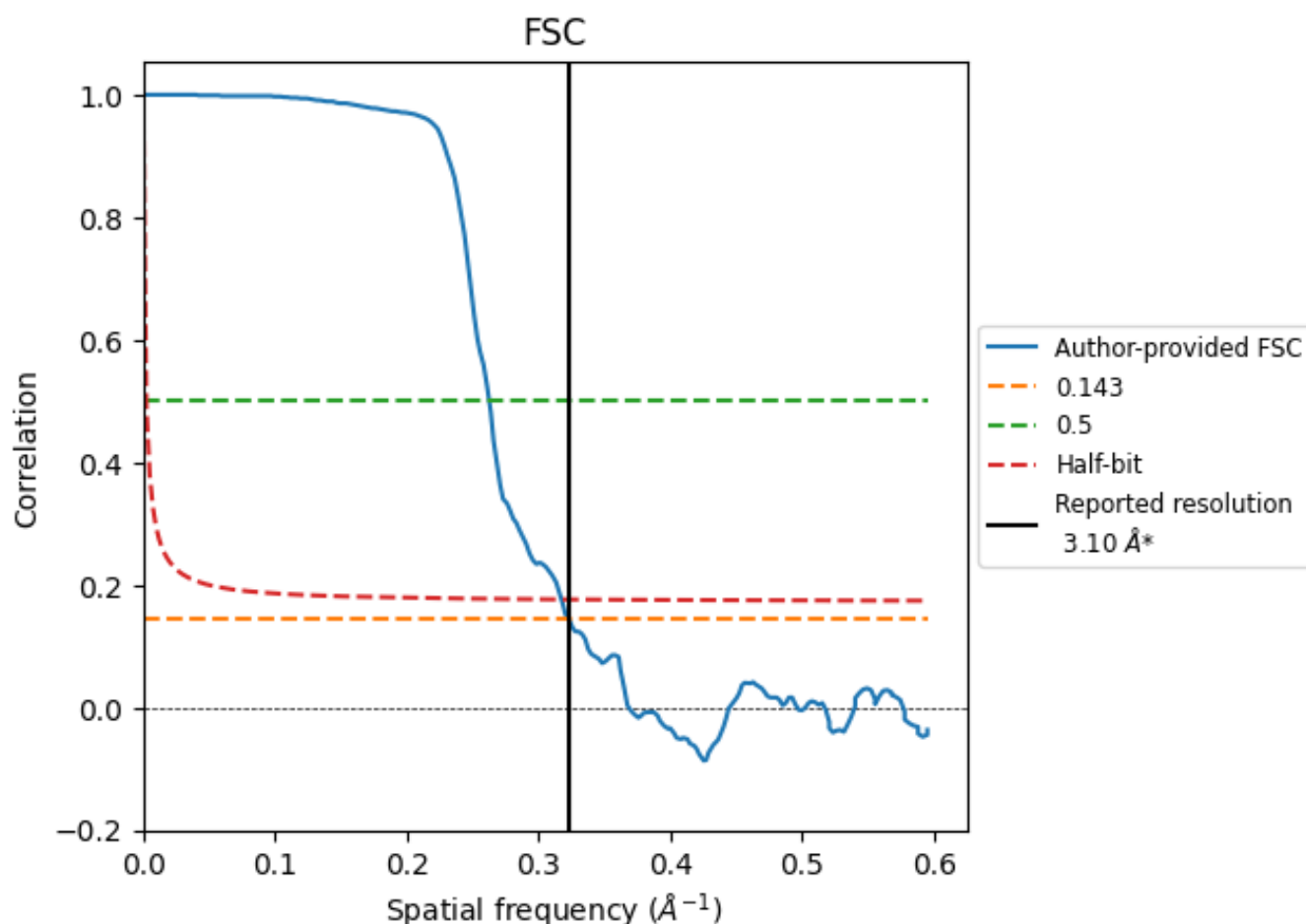


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 0.323 Å⁻¹

8.2 Resolution estimates [i](#)

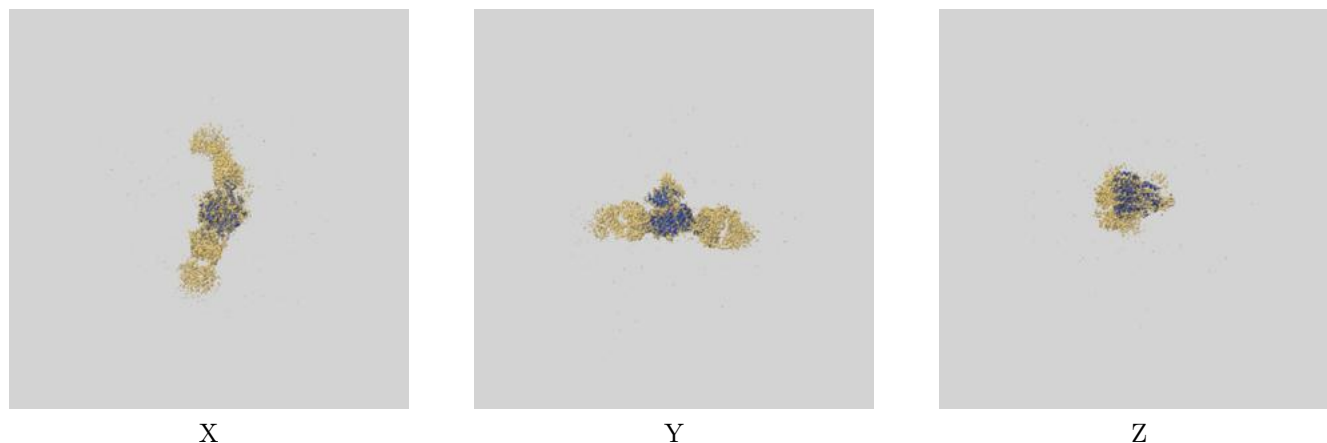
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.09	3.81	3.15
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

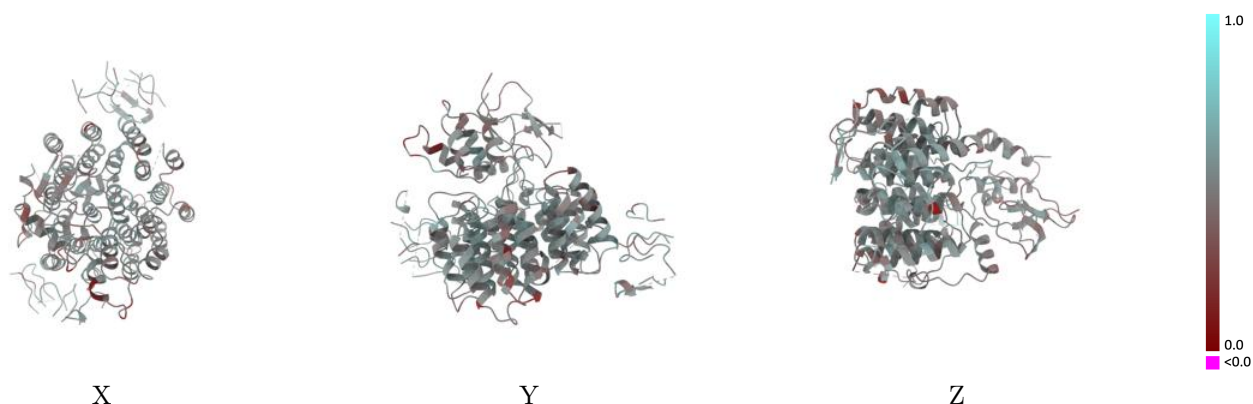
This section contains information regarding the fit between EMDB map EMD-24743 and PDB model 7UIH. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



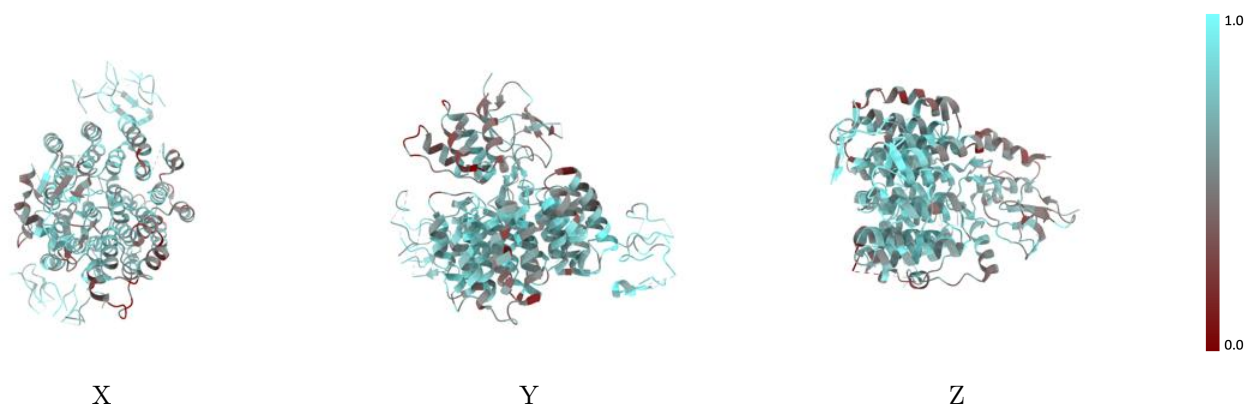
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



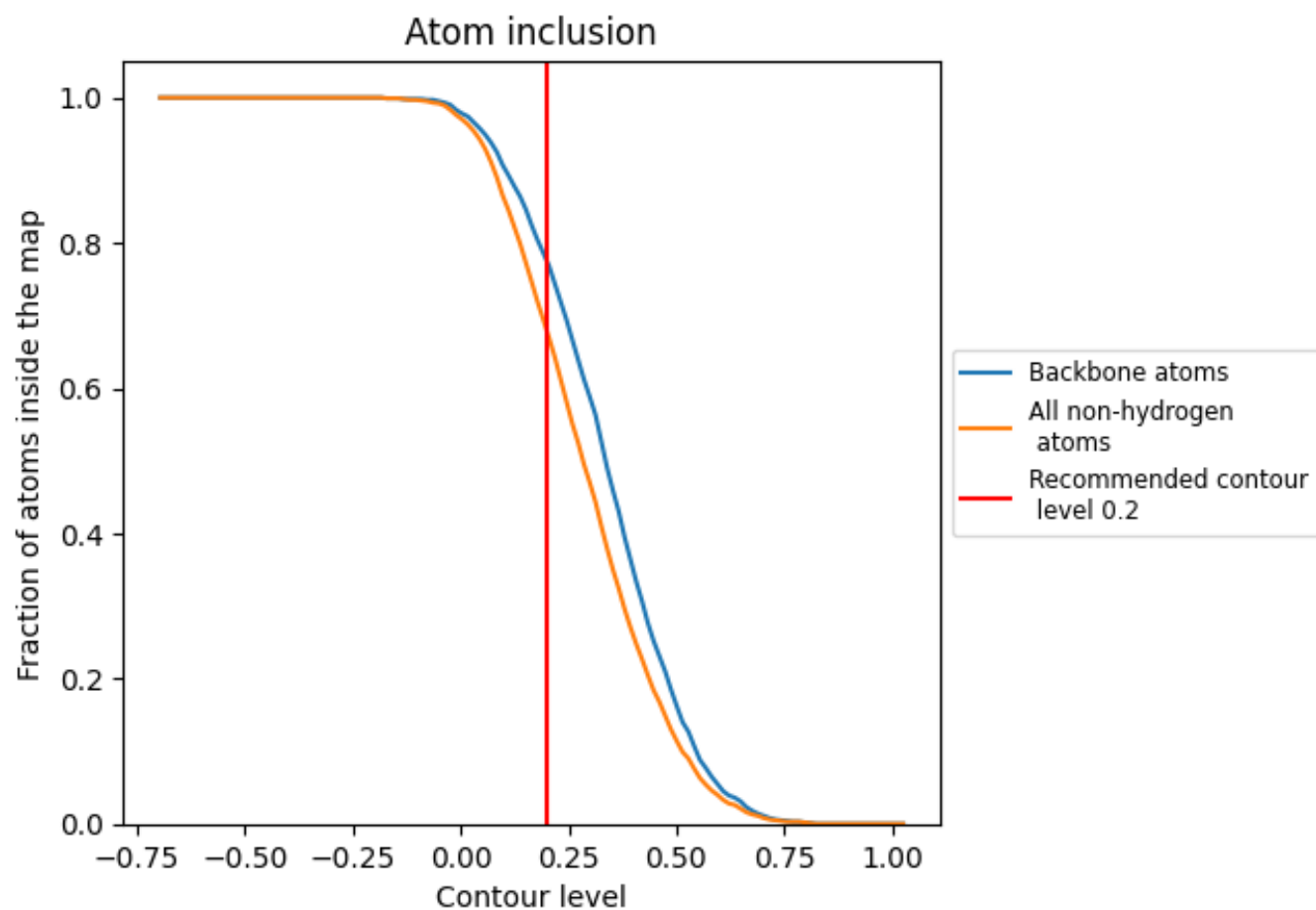
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 68% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6790	0.4910
A	0.6600	0.4860
C	0.7840	0.5150
D	0.8070	0.5100
E	0.7770	0.5250
F	0.8010	0.5220

