



wwPDB EM Validation Summary Report ⓘ

Jun 1, 2024 – 12:34 PM EDT

PDB ID : 7UIG
EMDB ID : EMD-26545
Title : Mediator-PIC Early (Mediator A)
Authors : Gorbea Colon, J.J.; Chen, S.-F.; Tsai, K.L.; Murakami, K.
Deposited on : 2022-03-29
Resolution : 4.30 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.dev92
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

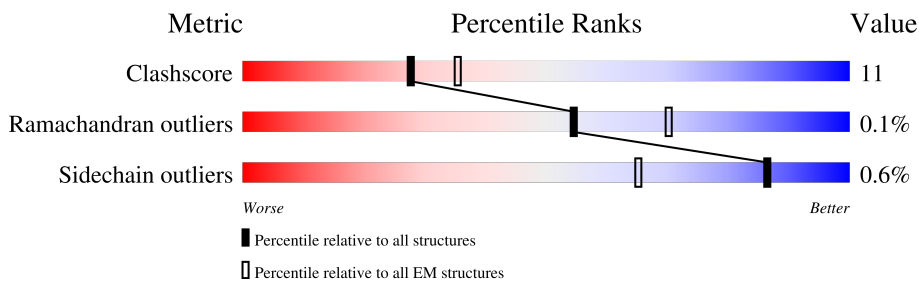
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.30 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	566	
2	d	284	
3	f	295	
4	g	222	
5	h	223	
6	i	149	
7	j	157	
8	k	115	

Continued on next page...

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Mol	Chain	Length	Quality of chain
9	n	1082	
10	q	687	
11	r	307	
12	s	220	
13	t	210	
14	u	140	
15	v	121	
16	w	127	
17	z	1733	

2 Entry composition [i](#)

There are 17 unique types of molecules in this entry. The entry contains 27580 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 Mediator of RNA polymerase II transcription subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	a	365	3008	1932	478	588	10	0	0

- Molecule 2 is a protein called Mediator of RNA polymerase II transcription subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	d	171	1388	875	233	276	4	0	0

- Molecule 3 is a protein called Mediator of RNA polymerase II transcription subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	f	169	1407	905	234	262	6	0	0

- Molecule 4 is a protein called Mediator of RNA polymerase II transcription subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	g	169	1409	903	238	263	5	0	0

- Molecule 5 is a protein called Mediator of RNA polymerase II transcription subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	h	136	1126	709	199	215	3	0	0

- Molecule 6 is a protein called Mediator of RNA polymerase II transcription subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	i	83	709	444	130	134	1	0	0

- Molecule 7 is a protein called Mediator of RNA polymerase II transcription subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	j	146	1173	725	206	239	3	0	0

- Molecule 8 is a protein called Mediator of RNA polymerase II transcription subunit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	k	108	876	546	149	177	4	0	0

- Molecule 9 is a protein called Mediator of RNA polymerase II transcription subunit 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	n	625	5139	3318	884	913	24	0	0

- Molecule 10 is a protein called Mediator of RNA polymerase II transcription subunit 17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	q	515	4182	2674	707	788	13	0	0

- Molecule 11 is a protein called Mediator of RNA polymerase II transcription subunit 18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	r	253	1995	1271	331	383	10	0	0

- Molecule 12 is a protein called Mediator of RNA polymerase II transcription subunit 19.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	s	81	657	415	109	132	1	0	0

- Molecule 13 is a protein called Mediator of RNA polymerase II transcription subunit 20.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	t	210	1609	1016	270	317	6	0	0

- Molecule 14 is a protein called Mediator of RNA polymerase II transcription subunit 21.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	u	122	Total	C	N	O	S	0	0
			978	611	163	199	5		

- Molecule 15 is a protein called Mediator of RNA polymerase II transcription subunit 22.

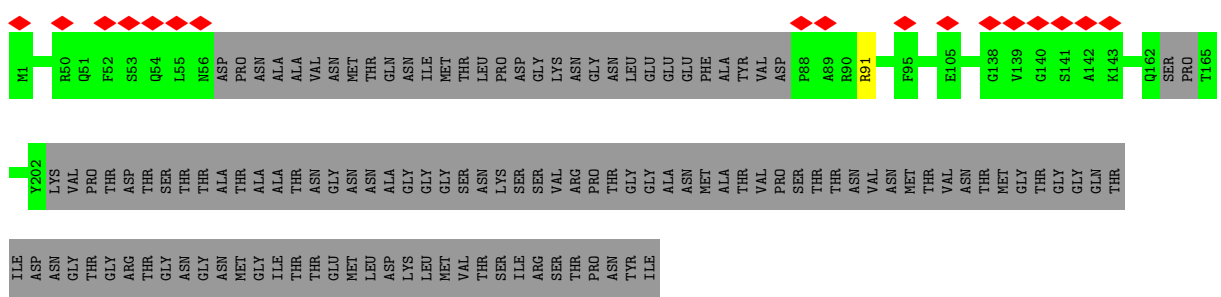
Mol	Chain	Residues	Atoms					AltConf	Trace
15	v	109	Total	C	N	O	S	0	0
			869	540	143	180	6		

- Molecule 16 is a protein called Mediator of RNA polymerase II transcription subunit 31.

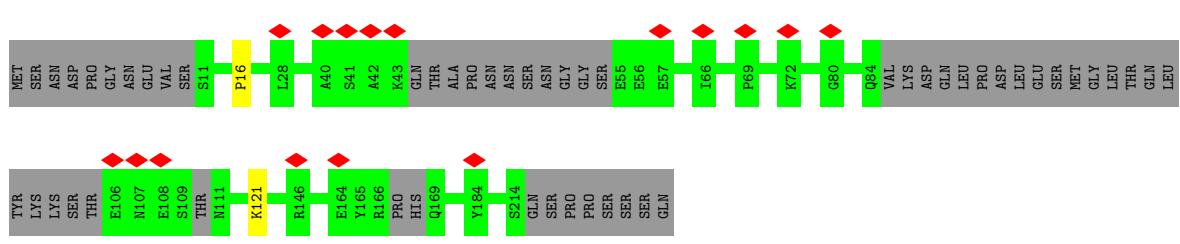
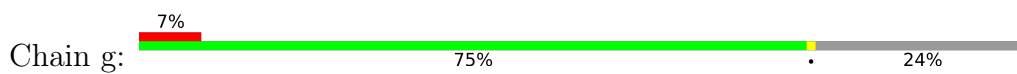
Mol	Chain	Residues	Atoms					AltConf	Trace
16	w	103	Total	C	N	O	S	0	0
			871	575	135	155	6		

- Molecule 17 is a protein called DNA-directed RNA polymerase II subunit RPB1.

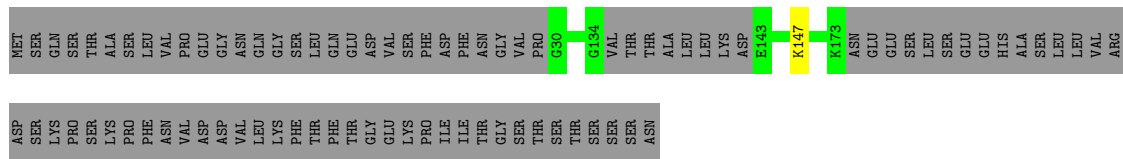
Mol	Chain	Residues	Atoms				AltConf	Trace
17	z	25	Total	C	N	O	0	0
			184	116	25	43		



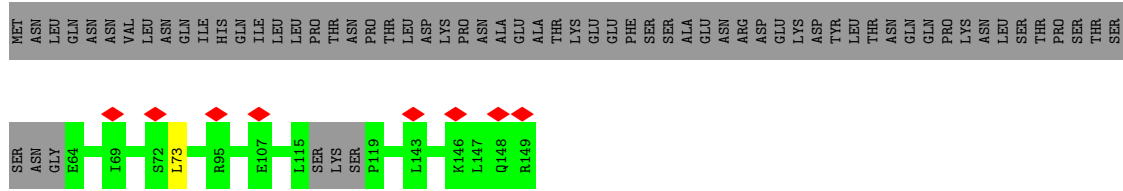
• Molecule 4: Mediator of RNA polymerase II transcription subunit 7



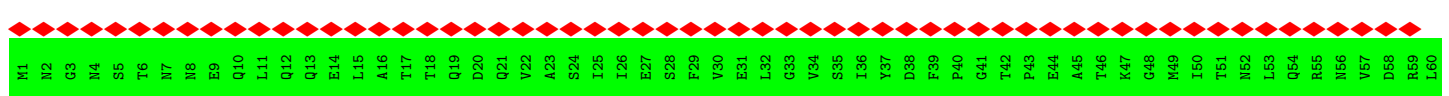
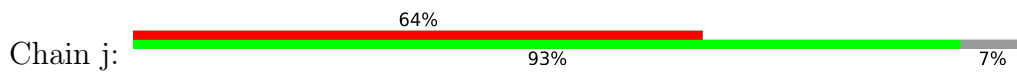
• Molecule 5: Mediator of RNA polymerase II transcription subunit 8



• Molecule 6: Mediator of RNA polymerase II transcription subunit 9



• Molecule 7: Mediator of RNA polymerase II transcription subunit 10



SER
PRO
SER
TYR
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THR
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SER
PRO
LYS
GLN
ASP
GLU
GLN
LYS
HIS
ASN
GLU
ASN
GLU
ASN
SER
ARG

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	341314	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	64000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.071	Depositor
Minimum map value	-0.016	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.02	Depositor
Map size (\AA)	496.8, 496.8, 496.8	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.38, 1.38, 1.38	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	0.65	0/3067	0.86	4/4148 (0.1%)
2	d	0.27	0/1405	0.58	1/1889 (0.1%)
3	f	0.27	0/1440	0.50	0/1946
4	g	0.28	0/1434	0.50	1/1930 (0.1%)
5	h	0.26	0/1147	0.55	0/1552
6	i	0.26	0/720	0.63	1/965 (0.1%)
7	j	0.24	0/1188	0.45	0/1604
8	k	0.27	0/885	0.49	0/1183
9	n	0.24	0/5226	0.46	0/7051
10	q	0.26	0/4245	0.49	0/5702
11	r	0.26	0/2030	0.50	0/2747
12	s	0.24	0/669	0.43	0/906
13	t	0.26	0/1635	0.50	0/2215
14	u	0.29	0/984	0.58	0/1317
15	v	0.26	0/873	0.51	0/1177
16	w	0.26	0/897	0.42	0/1219
17	z	0.25	0/194	0.41	0/270
All	All	0.33	0/28039	0.55	7/37821 (0.0%)

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	a	349	ARG	NE-CZ-NH1	7.31	123.95	120.30
1	a	67	TYR	CB-CG-CD2	-6.32	117.21	121.00
1	a	189	ARG	NE-CZ-NH1	6.31	123.46	120.30
6	i	73	LEU	CA-CB-CG	5.86	128.77	115.30
1	a	311	ARG	NE-CZ-NH1	5.41	123.01	120.30

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	3008	0	2953	0	0
2	d	1388	0	1400	0	0
3	f	1407	0	1385	0	0
4	g	1409	0	1434	0	0
5	h	1126	0	1125	0	0
6	i	709	0	712	0	0
7	j	1173	0	1156	0	0
8	k	876	0	885	0	0
9	n	5139	0	5375	0	0
10	q	4182	0	4324	0	0
11	r	1995	0	2018	0	0
12	s	657	0	636	0	0
13	t	1609	0	1626	0	0
14	u	978	0	1005	0	0
15	v	869	0	881	0	0
16	w	871	0	857	0	0
17	z	184	0	163	0	0
All	All	27580	0	27935	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 11.

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	357/566 (63%)	343 (96%)	11 (3%)	3 (1%)	19	60
2	d	165/284 (58%)	164 (99%)	1 (1%)	0	100	100
3	f	163/295 (55%)	161 (99%)	2 (1%)	0	100	100
4	g	159/222 (72%)	158 (99%)	1 (1%)	0	100	100
5	h	132/223 (59%)	129 (98%)	3 (2%)	0	100	100
6	i	79/149 (53%)	79 (100%)	0	0	100	100
7	j	142/157 (90%)	137 (96%)	5 (4%)	0	100	100
8	k	104/115 (90%)	104 (100%)	0	0	100	100
9	n	611/1082 (56%)	607 (99%)	4 (1%)	0	100	100
10	q	505/687 (74%)	500 (99%)	5 (1%)	0	100	100
11	r	249/307 (81%)	244 (98%)	5 (2%)	0	100	100
12	s	77/220 (35%)	76 (99%)	1 (1%)	0	100	100
13	t	208/210 (99%)	206 (99%)	2 (1%)	0	100	100
14	u	116/140 (83%)	115 (99%)	1 (1%)	0	100	100
15	v	105/121 (87%)	105 (100%)	0	0	100	100
16	w	99/127 (78%)	97 (98%)	2 (2%)	0	100	100
17	z	23/1733 (1%)	21 (91%)	1 (4%)	1 (4%)	2	25
All	All	3294/6638 (50%)	3246 (98%)	44 (1%)	4 (0%)	54	85

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
17	z	21	SER
1	a	291	SER
1	a	306	ASN
1	a	289	CYS

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	350/528 (66%)	347 (99%)	3 (1%)	78	88
2	d	158/258 (61%)	157 (99%)	1 (1%)	86	92
3	f	158/259 (61%)	157 (99%)	1 (1%)	86	92
4	g	160/208 (77%)	159 (99%)	1 (1%)	86	92
5	h	128/207 (62%)	127 (99%)	1 (1%)	81	89
6	i	82/144 (57%)	82 (100%)	0	100	100
7	j	134/145 (92%)	134 (100%)	0	100	100
8	k	101/108 (94%)	101 (100%)	0	100	100
9	n	591/1001 (59%)	588 (100%)	3 (0%)	88	93
10	q	482/642 (75%)	480 (100%)	2 (0%)	91	94
11	r	228/280 (81%)	228 (100%)	0	100	100
12	s	75/195 (38%)	75 (100%)	0	100	100
13	t	178/178 (100%)	178 (100%)	0	100	100
14	u	115/132 (87%)	110 (96%)	5 (4%)	29	55
15	v	101/113 (89%)	100 (99%)	1 (1%)	76	86
16	w	97/117 (83%)	97 (100%)	0	100	100
17	z	25/1520 (2%)	25 (100%)	0	100	100
All	All	3163/6035 (52%)	3145 (99%)	18 (1%)	86	92

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
14	u	100	LYS
15	v	45	ASN
14	u	137	LYS
9	n	449	ARG
14	u	4	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
14	u	9	GLN
14	u	98	GLN
5	h	150	GLN
8	k	19	GLN
9	n	715	ASN

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 monosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
10	q	1
14	u	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	q	318:ASN	C	319:LYS	N	5.82
1	u	80:LEU	C	81:PRO	N	3.18

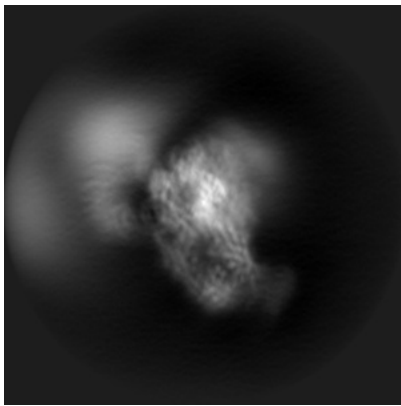
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26545. These allow visual inspection of the internal detail of the map and identification of artifacts.

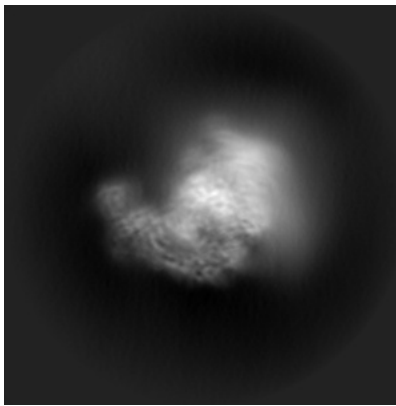
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

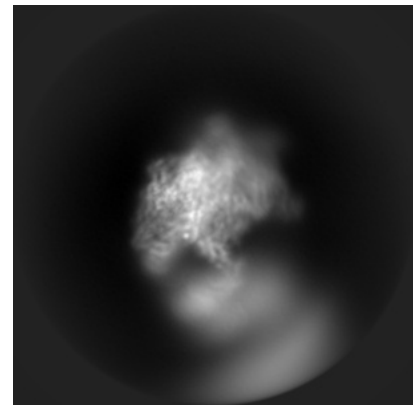
6.1.1 Primary map



X

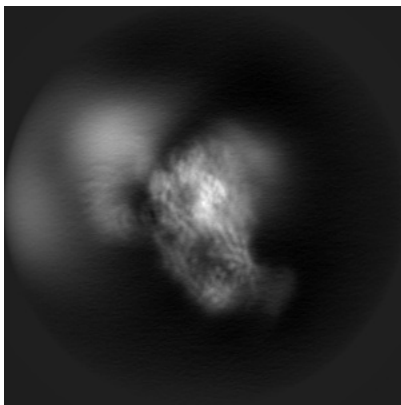


Y

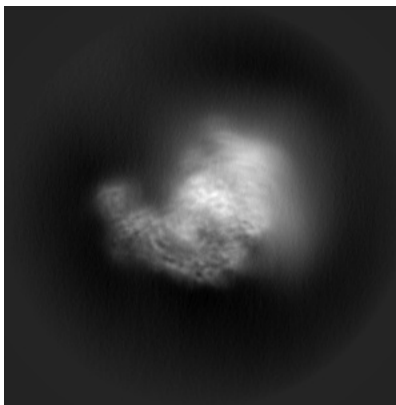


Z

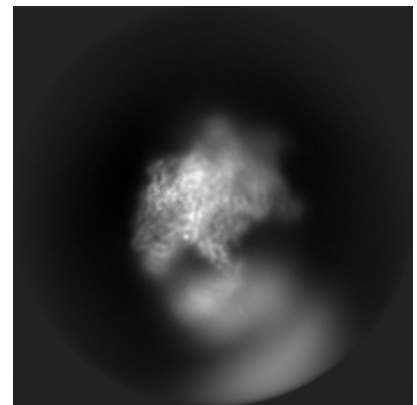
6.1.2 Raw map



X



Y

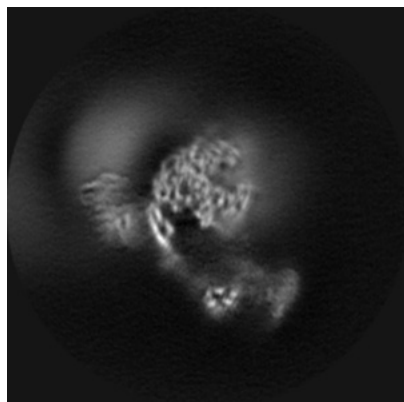


Z

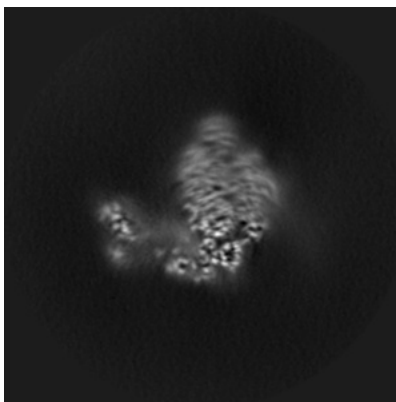
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

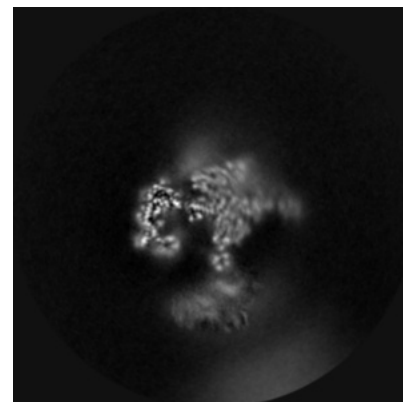
6.2.1 Primary map



X Index: 180

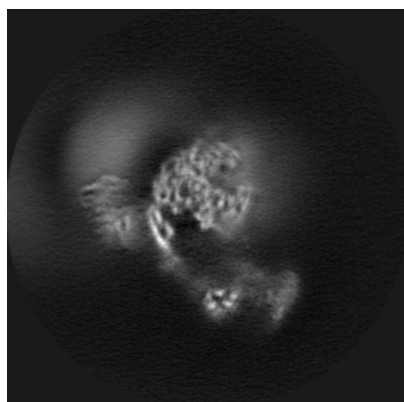


Y Index: 180

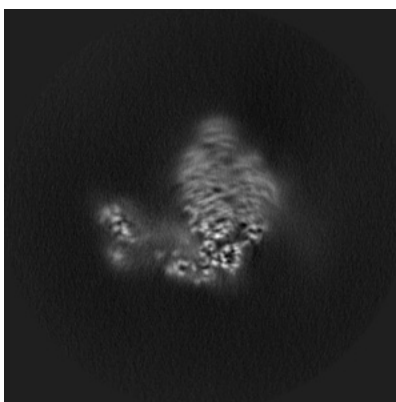


Z Index: 180

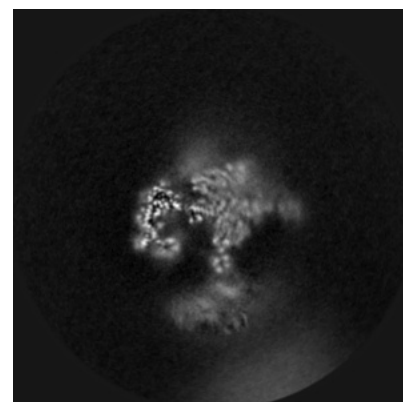
6.2.2 Raw map



X Index: 180



Y Index: 180

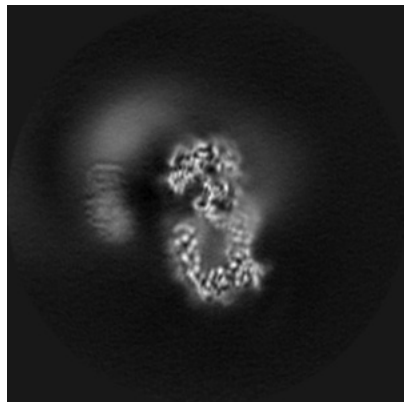


Z Index: 180

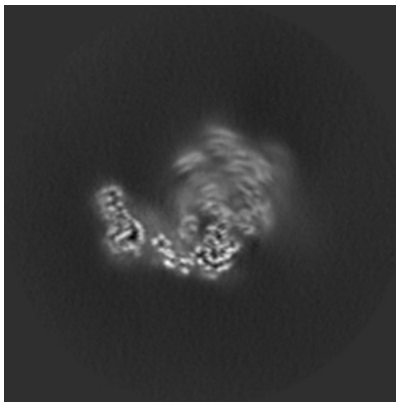
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

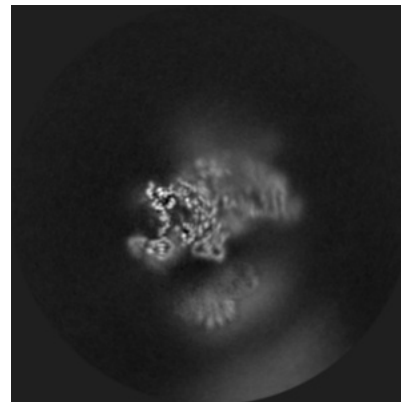
6.3.1 Primary map



X Index: 161

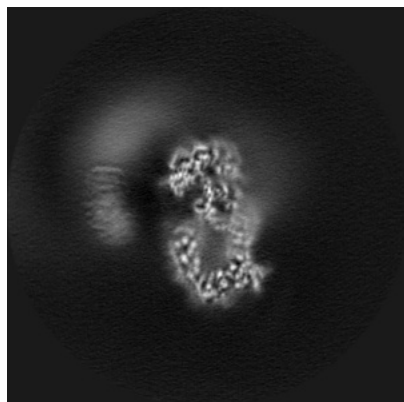


Y Index: 192

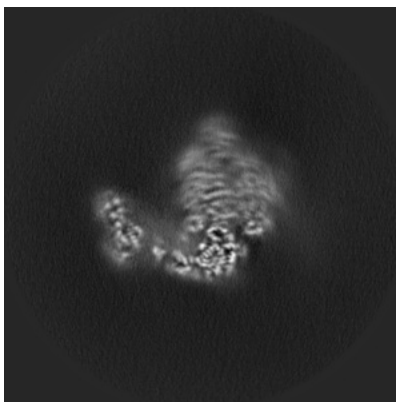


Z Index: 199

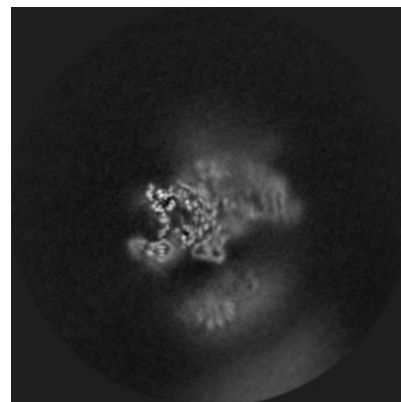
6.3.2 Raw map



X Index: 162



Y Index: 185

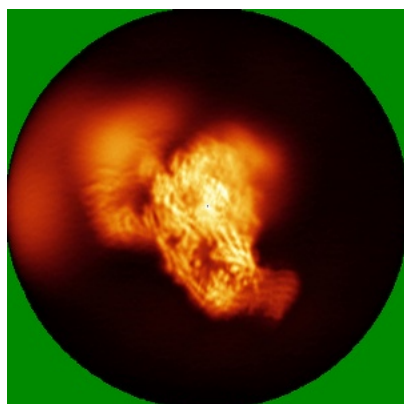


Z Index: 199

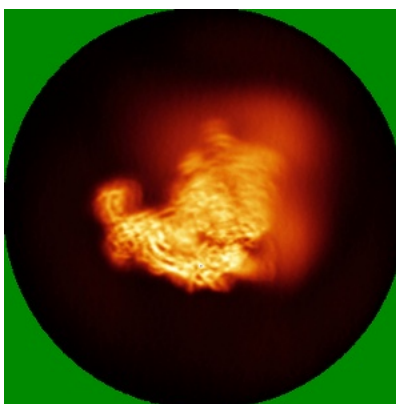
The images above show the largest variance slices of the map in three orthogonal directions.

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

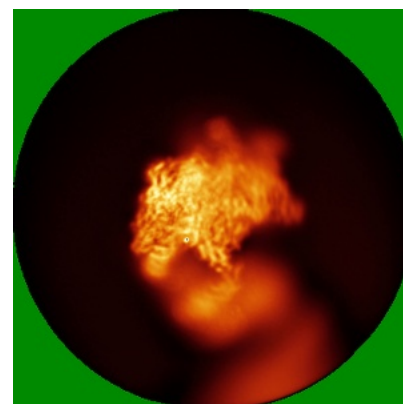
6.4.1 Primary map



X

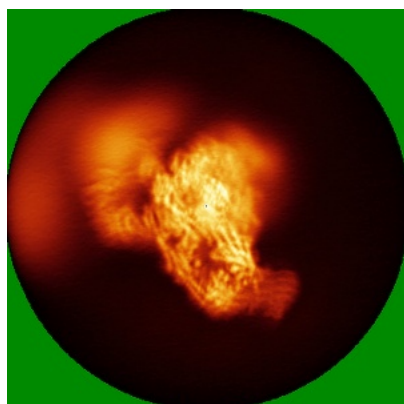


Y

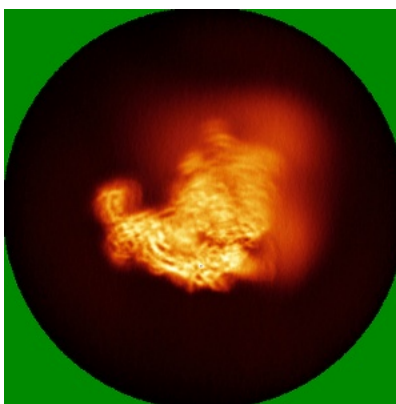


Z

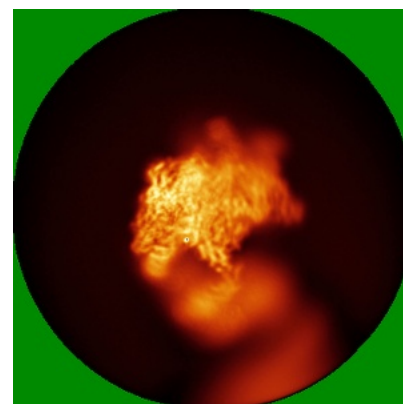
6.4.2 Raw map



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.

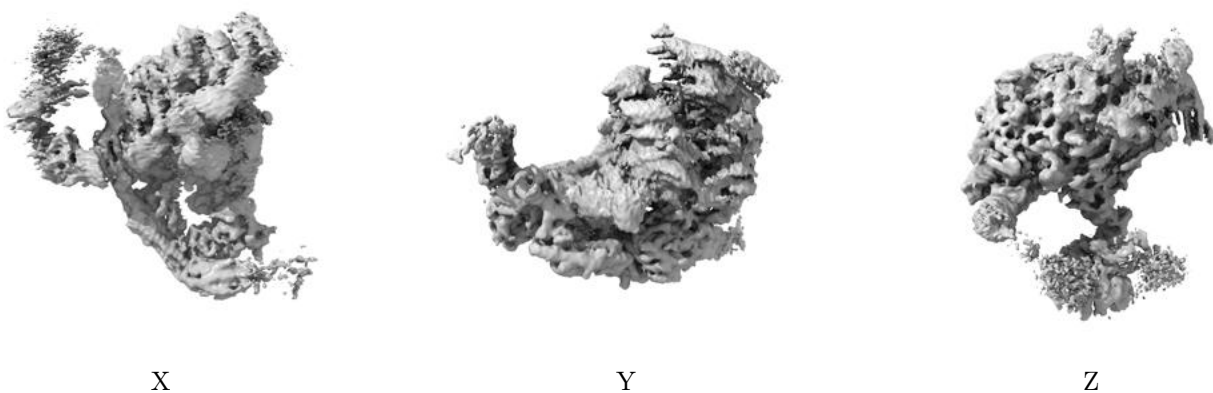
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.02. 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.

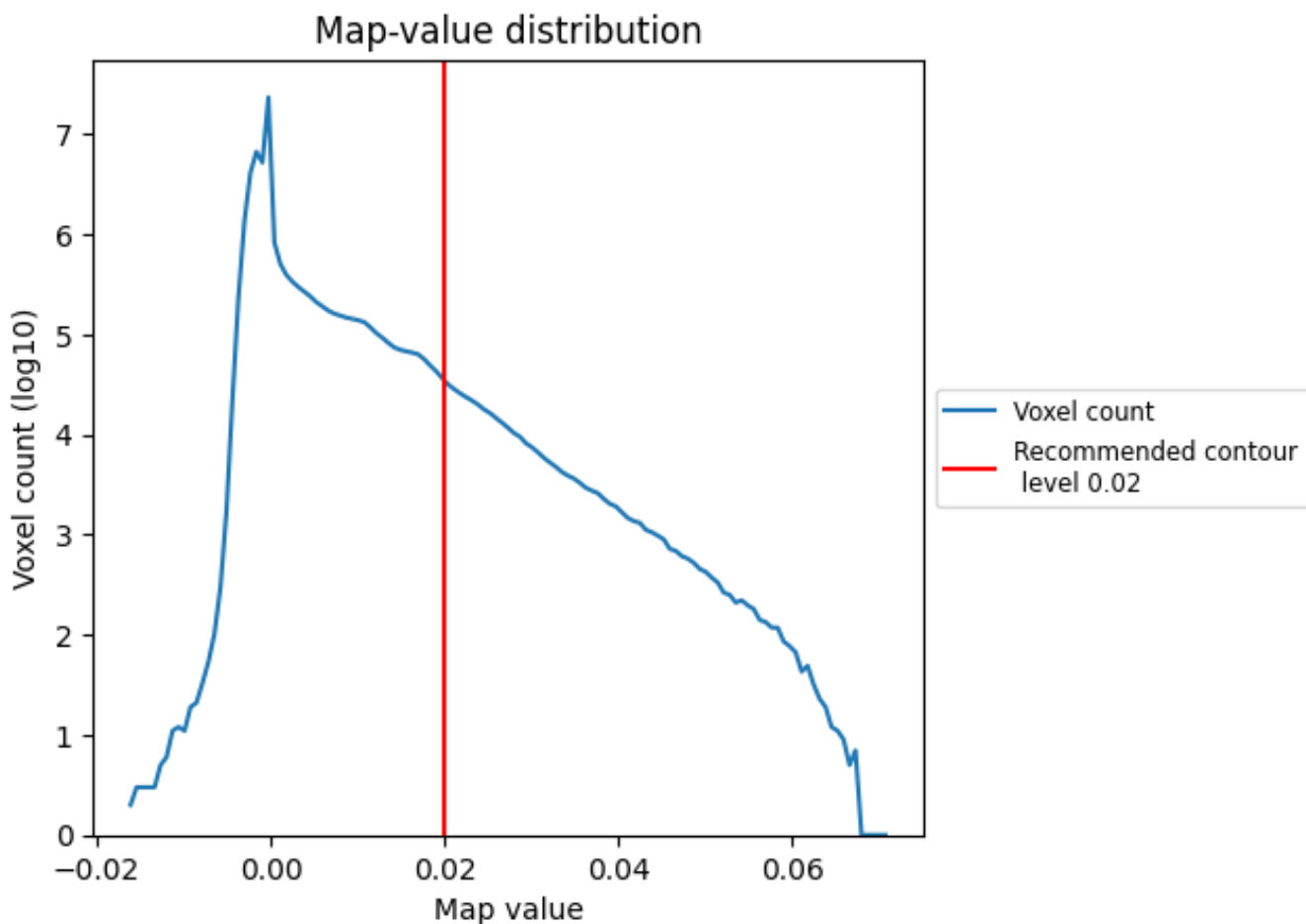
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

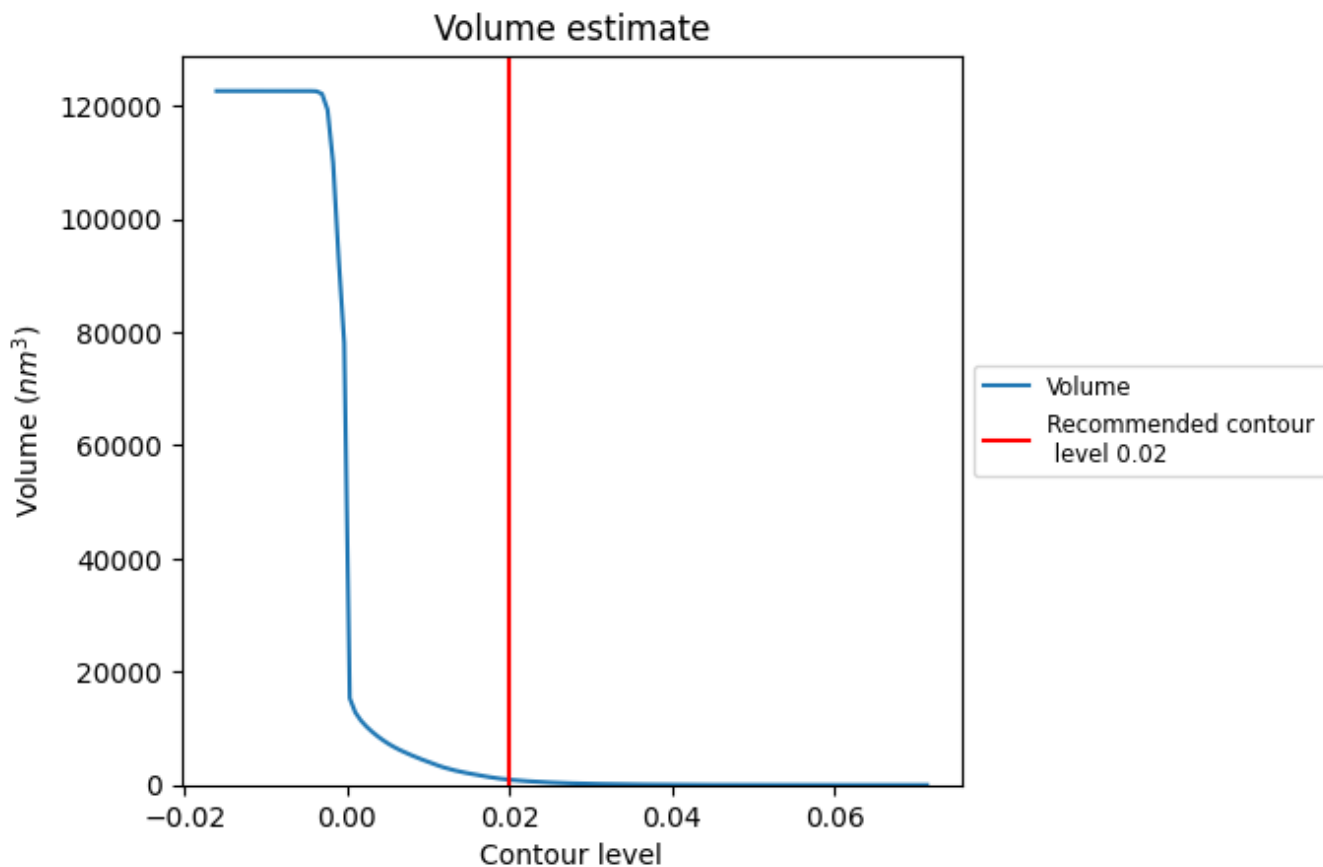
This section contains the results of statistical analysis of the map.

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.

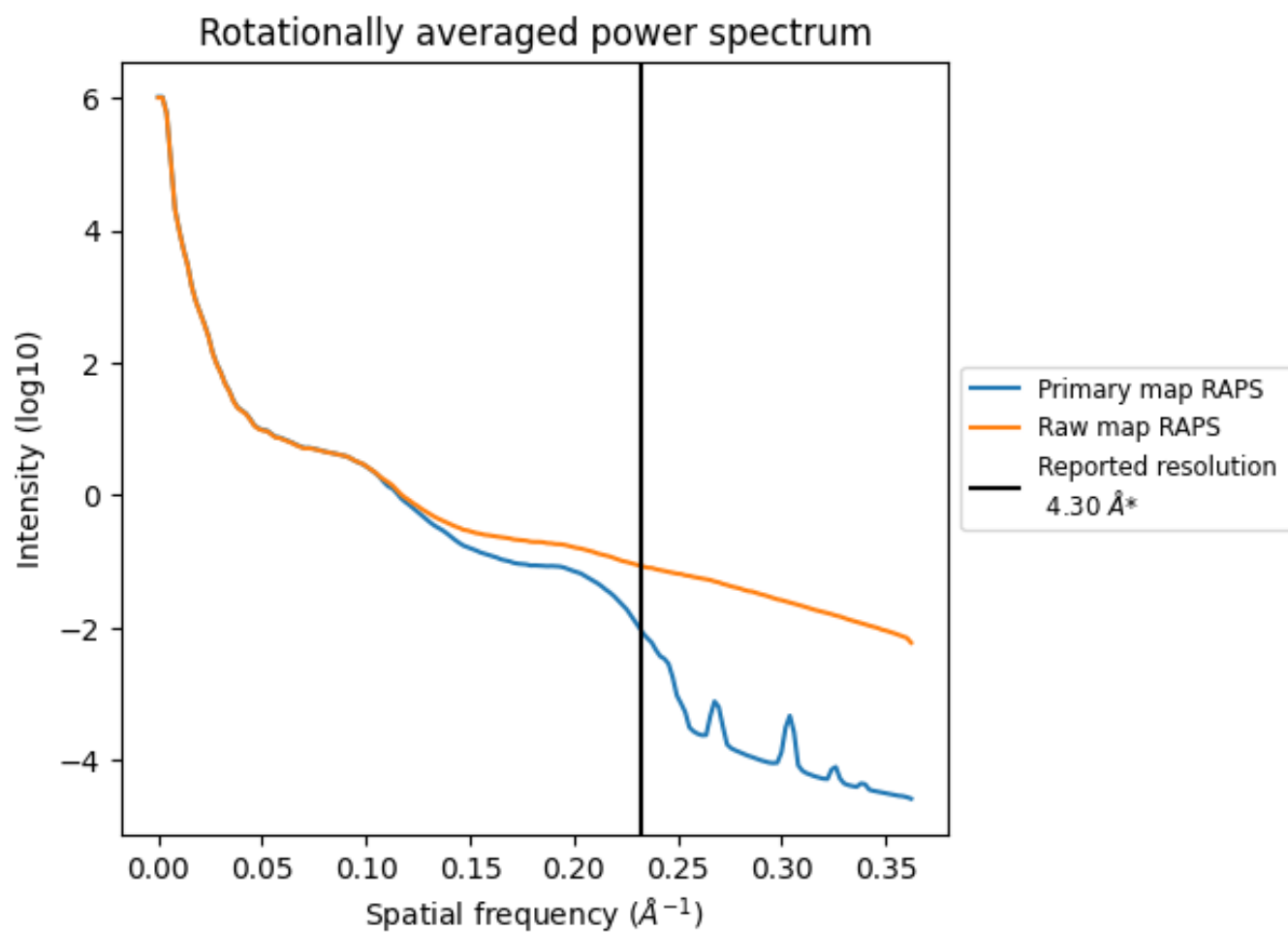
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 925 nm³; this corresponds to an approximate mass of 836 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](#)

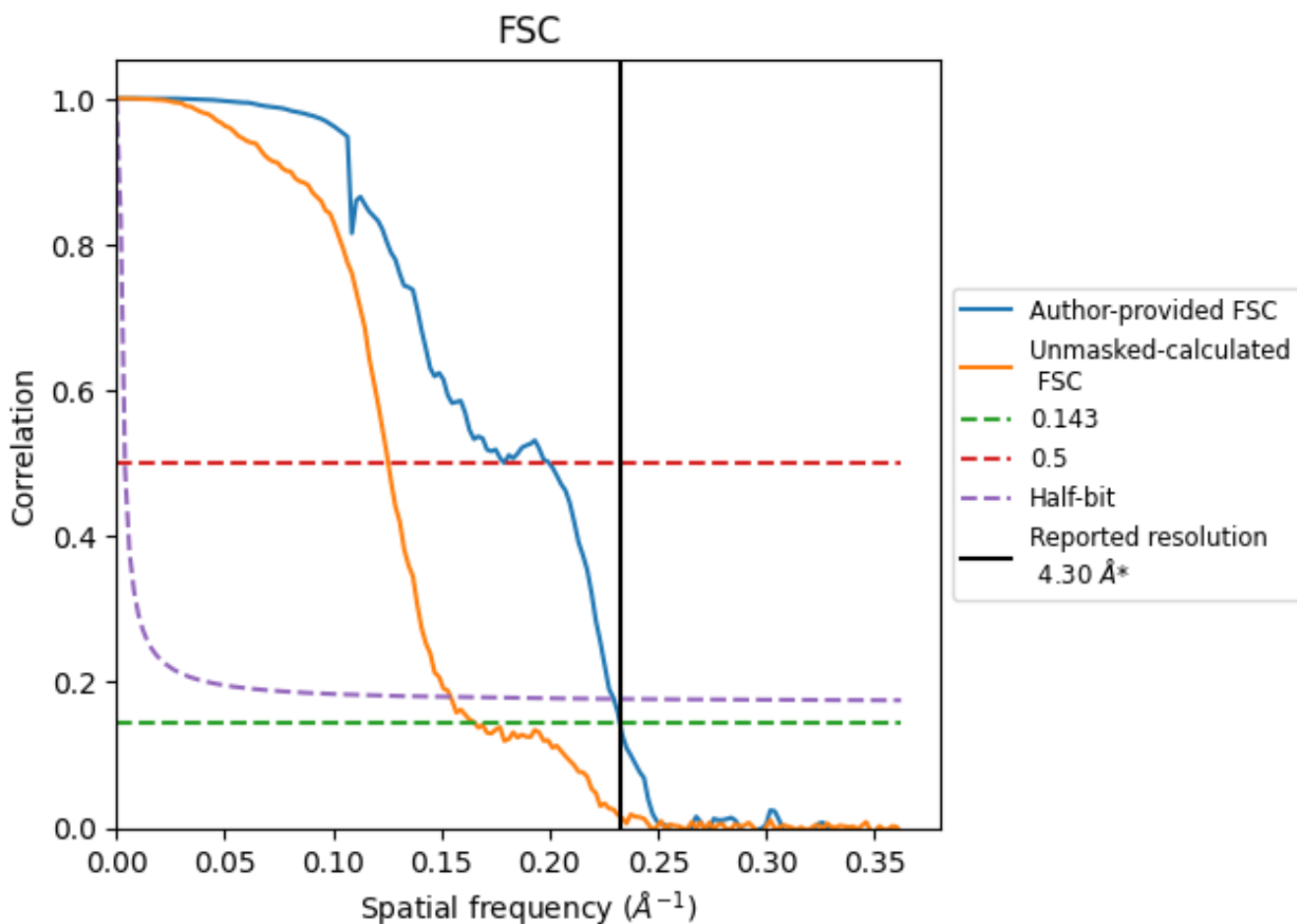


*Reported resolution corresponds to spatial frequency of 0.233 Å⁻¹

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 0.233 Å⁻¹

8.2 Resolution estimates [i](#)

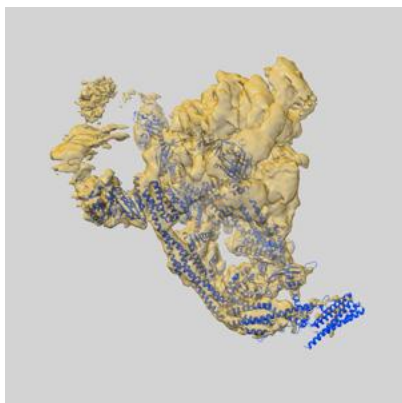
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.30	-	-
Author-provided FSC curve	4.30	5.01	4.36
Unmasked-calculated*	6.04	7.97	6.48

*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 6.04 differs from the reported value 4.3 by more than 10 %

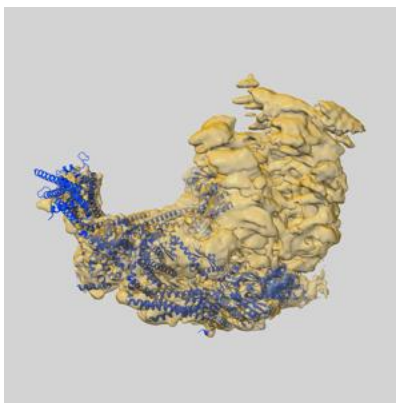
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26545 and PDB model 7UIG. Per-residue inclusion information can be found in section 3 on page 7.

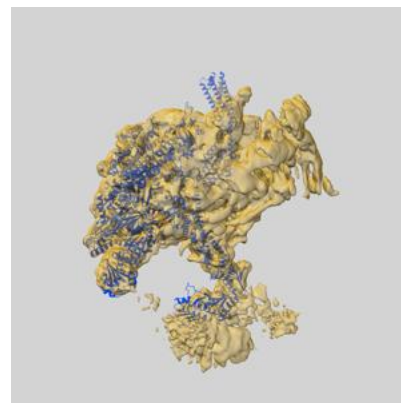
9.1 Map-model overlay [i](#)



X



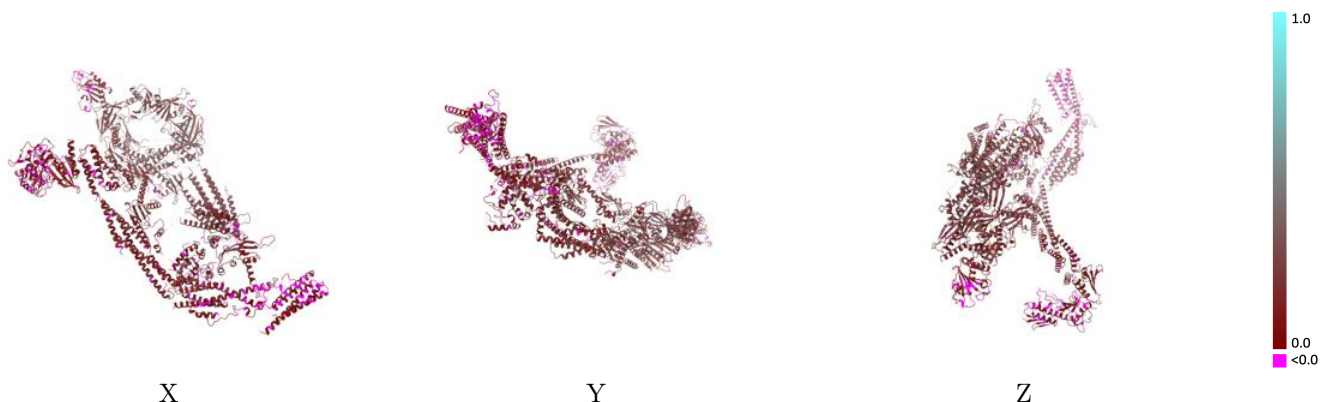
Y



Z

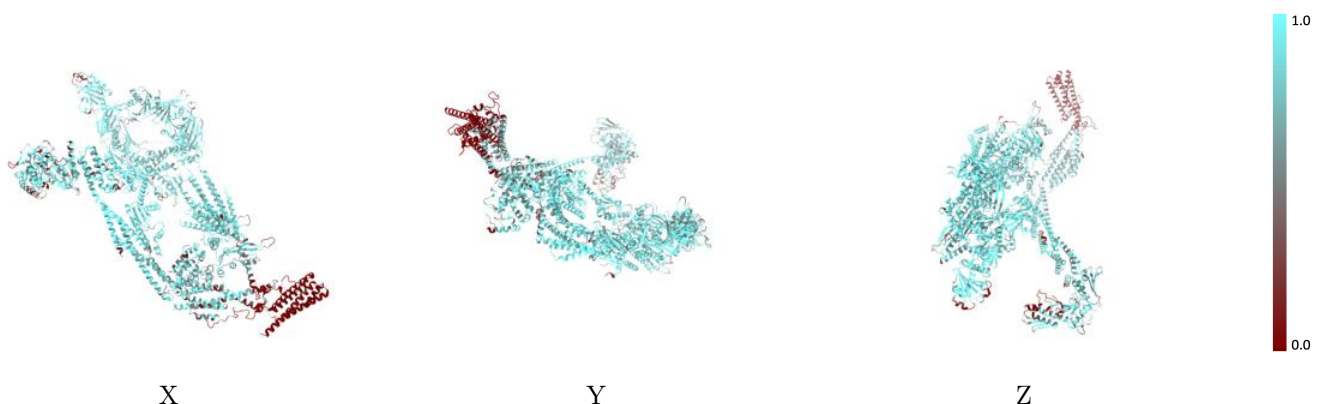
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



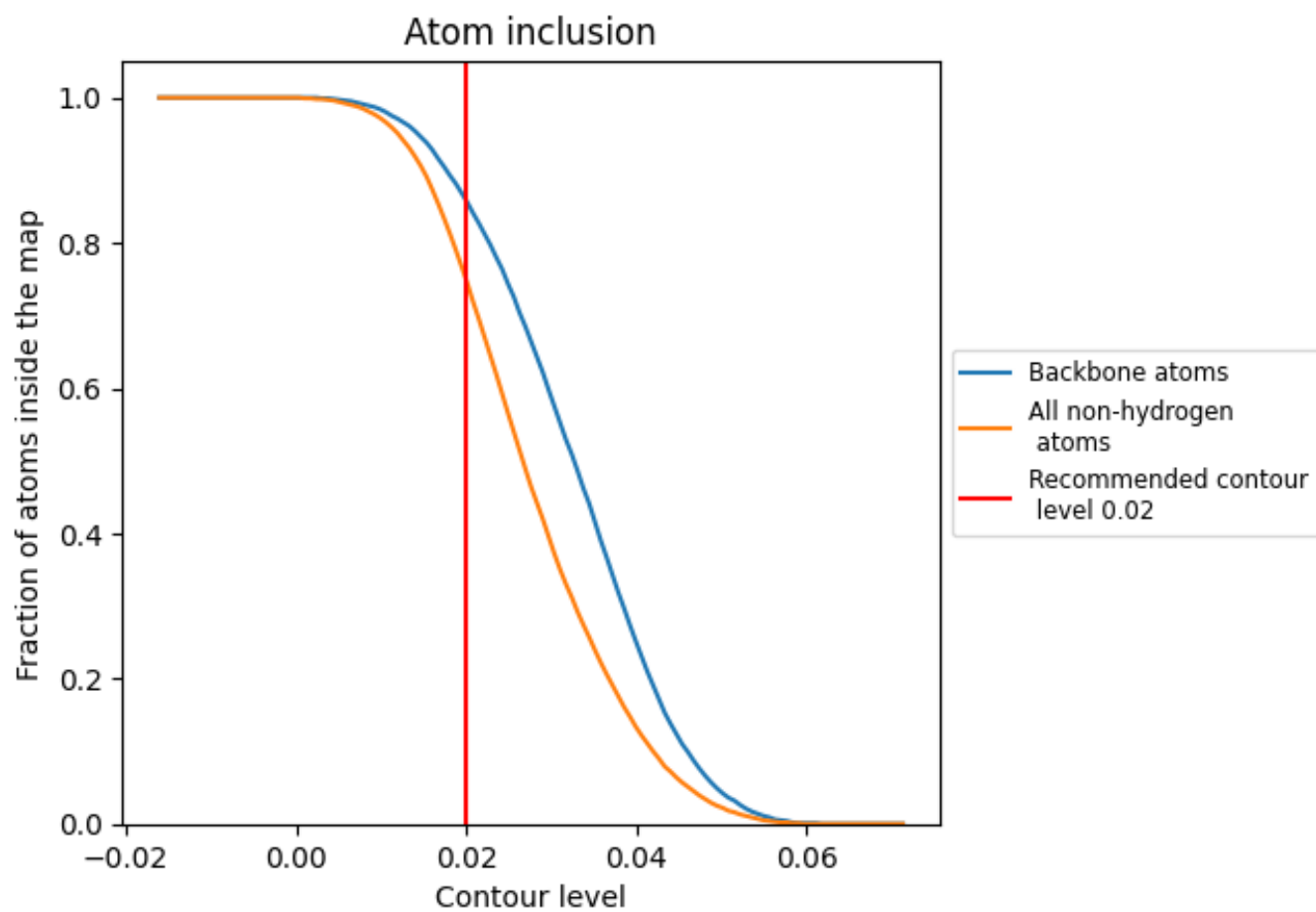
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).





































9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 75% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7500	 0.1720
a	 0.6590	 0.0950
d	 0.7950	 0.1550
f	 0.7620	 0.1840
g	 0.7640	 0.1510
h	 0.9020	 0.1990
i	 0.8010	 0.1410
j	 0.2670	 0.0730
k	 0.8860	 0.2220
n	 0.7550	 0.1580
q	 0.8280	 0.2030
r	 0.8430	 0.2790
s	 0.1060	 0.0410
t	 0.8230	 0.2560
u	 0.7620	 0.1470
v	 0.8510	 0.1980
w	 0.8150	 0.1720
z	 0.8560	 0.1600

