



wwPDB EM Validation Summary Report ⓘ

May 19, 2025 – 08:58 PM EDT

PDB ID : 8VG1 / pdb_00008vg1
EMDB ID : EMD-43197
Title : Cryo-EM structure of FoxA1 and GATA4 in complex with ALBN1 nucleosome
Authors : Zhou, B.R.; Bai, Y.
Deposited on : 2023-12-22
Resolution : 2.48 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.dev118
MolProbity : 4-5-2 with Phenix2.0rc1
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

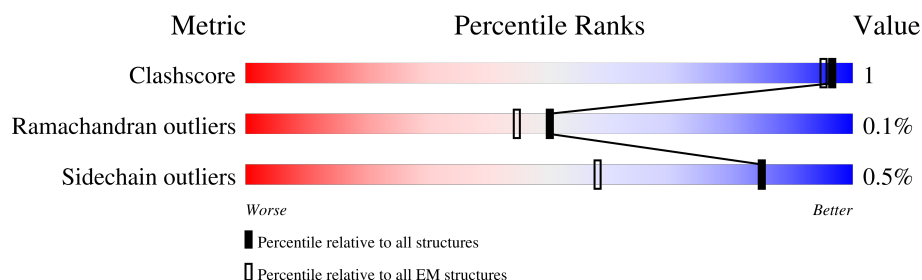
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.48 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	136	69% 29%
1	E	136	71% 29%
2	B	103	81% 19%
2	F	103	80% 19%
3	C	130	84% 16%
3	G	130	88% 11%
4	D	126	74% 25%
4	H	126	73% 25%

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Mol	Chain	Length	Quality of chain
5	I	186	<div><div></div><div>87%</div><div>5%8%</div></div>
6	J	186	<div><div></div><div>91%</div><div>8%</div></div>
7	O	478	<div><div>6%19%</div><div></div><div>79%</div></div>
8	T	835	<div><div>6%</div><div></div><div>94%</div></div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14434 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 Histone H3.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	97	Total	C	N	O	S	0	0
			799	505	153	137	4		
1	E	97	Total	C	N	O	S	0	0
			799	505	153	137	4		

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	83	Total	C	N	O	S	0	0
			662	418	129	114	1		
2	F	83	Total	C	N	O	S	0	0
			662	418	129	114	1		

- Molecule 3 is a protein called Histone H2A type 1-B/E.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	109	Total	C	N	O	0	0
			840	529	166	145		
3	G	116	Total	C	N	O	0	0
			900	565	179	156		

- Molecule 4 is a protein called Histone H2B type 1-J.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	95	Total	C	N	O	S	0	0
			745	468	136	139	2		
4	H	95	Total	C	N	O	S	0	0
			745	468	136	139	2		

- Molecule 5 is a DNA chain called DNA (171-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	171	Total	C	N	O	P	0	0
			3501	1673	619	1038	171		

- Molecule 6 is a DNA chain called DNA (171-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	171	Total	C	N	O	P	0	0
			3510	1672	653	1014	171		

- Molecule 7 is a protein called Hepatocyte nuclear factor 3-alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	O	102	Total	C	N	O	S	0	0
			856	548	152	149	7		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	473	HIS	-	expression tag	UNP P55317
O	474	HIS	-	expression tag	UNP P55317
O	475	HIS	-	expression tag	UNP P55317
O	476	HIS	-	expression tag	UNP P55317
O	477	HIS	-	expression tag	UNP P55317
O	478	HIS	-	expression tag	UNP P55317

- Molecule 8 is a protein called Maltose/maltodextrin-binding periplasmic protein, Transcriptin on factor GATA-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	T	53	Total	C	N	O	S	0	0
			414	253	83	72	6		

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
T	-392	MET	-	initiating methionine	UNP P0AEX9
T	-391	GLY	-	expression tag	UNP P0AEX9
T	-390	SER	-	expression tag	UNP P0AEX9
T	-389	SER	-	expression tag	UNP P0AEX9
T	-388	HIS	-	expression tag	UNP P0AEX9
T	-387	HIS	-	expression tag	UNP P0AEX9
T	-386	HIS	-	expression tag	UNP P0AEX9
T	-385	HIS	-	expression tag	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
T	-384	HIS	-	expression tag	UNP P0AEX9
T	-383	HIS	-	expression tag	UNP P0AEX9
T	-382	GLY	-	expression tag	UNP P0AEX9
T	-381	SER	-	expression tag	UNP P0AEX9
T	-380	SER	-	expression tag	UNP P0AEX9
T	-379	MET	-	expression tag	UNP P0AEX9
T	-12	ASN	-	linker	UNP P0AEX9
T	-11	GLY	-	linker	UNP P0AEX9
T	-10	ILE	-	linker	UNP P0AEX9
T	-9	GLU	-	linker	UNP P0AEX9
T	-8	GLU	-	linker	UNP P0AEX9
T	-7	ASN	-	linker	UNP P0AEX9
T	-6	LEU	-	linker	UNP P0AEX9
T	-5	TYR	-	linker	UNP P0AEX9
T	-4	PHE	-	linker	UNP P0AEX9
T	-3	GLN	-	linker	UNP P0AEX9
T	-2	SER	-	linker	UNP P0AEX9
T	-1	ASN	-	linker	UNP P0AEX9
T	0	ALA	-	linker	UNP P0AEX9

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

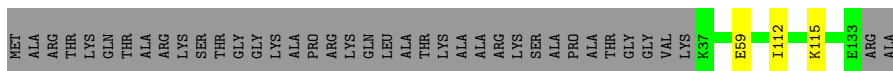
Mol	Chain	Residues	Atoms		AltConf
9	T	1	Total	Zn	0
			1	1	

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.

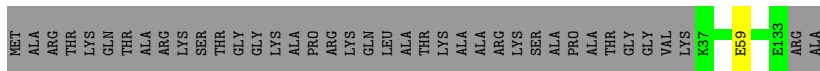
- Molecule 1: Histone H3.1

Chain A: 




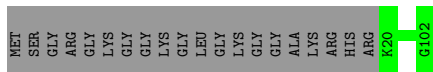
- Molecule 1: Histone H3.1

Chain E: 




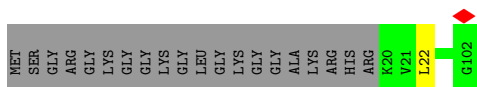
- Molecule 2: Histone H4

Chain B: 




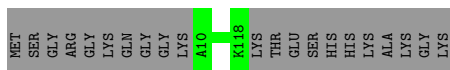
- Molecule 2: Histone H4

Chain F: 



- Molecule 3: Histone H2A type 1-B/E

Chain C: 



- Molecule 3: Histone H2A type 1-B/E

MET
 SER
 GLY
 ARG
 GLY
 LYS
 GLN
 GLY
 GLY
 LYS
 A10
 V114
 E121
 K125
 ALA
 LYS
 GLY
 LYS

- Chain D:  74% . 25%

MET	PRO	GLU	PRO	ALA	LYS	SER	ALA	PRO	PRO	LYS	LYS	GLY	SER	LYS	LYS	ALA	VAL	THR	LYS	ALA	GLN	LYS	ASP	GLY	LYS	LYS	ARG	K30	A107	V111	A124	LYS
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- Chain H:  73% 25%

NET	PRO	GLU	PRO	ALA	LYS	SER	ALA	PRO	ALA	LYS	GLY	SER	LYS	ALA	VAL	THR	LYS	ALA	GLN	LYS	LYS	ASP	GLY	LYS	LYS	ARG	K30	T90	E93	E113	A124	LYS
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	------	------	-----

- Chain I: 87% 5% 8%

DA DT DC DG DA DG DT DG DG DG DT DA DC DT T16 T19 G20 T21 G22 T23 G55 G68 T69 T89 T90 A120 T121 T167 T168 T186

- Chain J: 91% 8%

- Chain O:  6% 19% 79%

[illegible][illegible]

	ALA	ALA	SER	MET	ASN	GLY	LEU	GLY	PRO	Tyr	ALA	ALA	ALA	ALA	ALA	ALA	ALA	ALA	ASP	Lys	THR	PHE	Lys	ARG	SER	Tyr	PRO	H168	A199	Q195	A196	P187	S188	K189	M190	Q198	F202
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L203	F204	R208	L222	V231	A232	R233	S234	P235	D236	K237	P238	G239	S242	Y243	W244	T245	P248	D249	N252	M253	P254	E255	N256	G257	C258	K264	K267	C268	F269	LYS	GLN	PRO	GLY	GLY	ALA	GLY	GLY	GLY	GLY	GLY	SER	SER	SER	GLY	GLY	GLY	ALA	LYS	GLY	GLY	PRO
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

SER
HIS
HIS
HIS
HIS
HIS
HIS

Chain T: 6% 94%

VAL
LEU
ALA
ASP
SER
HIS
GLY
ASP
ILE
ILE
THR
ALA

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	403489	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.063	Depositor
Minimum map value	-0.022	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.004	Depositor
Map size (\AA)	316.8, 316.8, 316.8	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.056, 1.056, 1.056	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.16	0/811	0.30	0/1088
1	E	0.16	0/811	0.31	0/1088
2	B	0.16	0/669	0.32	0/894
2	F	0.16	0/669	0.32	0/894
3	C	0.16	0/850	0.28	0/1146
3	G	0.16	0/912	0.29	0/1228
4	D	0.16	0/756	0.29	0/1015
4	H	0.16	0/756	0.27	0/1015
5	I	0.25	0/3922	0.52	0/6052
6	J	0.25	0/3942	0.48	0/6081
7	O	0.15	0/883	0.36	0/1192
8	T	0.15	0/420	0.33	0/564
All	All	0.21	0/15401	0.42	0/22257

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	799	0	838	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	799	0	838	1	0
2	B	662	0	709	0	0
2	F	662	0	709	0	0
3	C	840	0	902	0	0
3	G	900	0	960	2	0
4	D	745	0	771	1	0
4	H	745	0	771	2	0
5	I	3501	0	1936	5	0
6	J	3510	0	1924	1	0
7	O	856	0	832	6	0
8	T	414	0	425	3	0
9	T	1	0	0	0	0
All	All	14434	0	11615	19	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 1.

The worst 5 of 19 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)
5:I:89:DT:H2''	5:I:90:DT:H72	1.83	0.60
7:O:222:LEU:HD11	7:O:244:TRP:CE2	2.39	0.57
4:H:113:GLU:OE2	8:T:320:ARG:NH1	2.37	0.56
7:O:253:MET:SD	7:O:253:MET:N	2.79	0.55
5:I:120:DA:H2''	5:I:121:DT:H72	1.92	0.50

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	95/136 (70%)	95 (100%)	0	0	100	100
1	E	95/136 (70%)	95 (100%)	0	0	100	100
2	B	81/103 (79%)	79 (98%)	2 (2%)	0	100	100
2	F	81/103 (79%)	80 (99%)	1 (1%)	0	100	100
3	C	107/130 (82%)	106 (99%)	1 (1%)	0	100	100
3	G	114/130 (88%)	113 (99%)	1 (1%)	0	100	100
4	D	93/126 (74%)	92 (99%)	1 (1%)	0	100	100
4	H	93/126 (74%)	92 (99%)	1 (1%)	0	100	100
7	O	100/478 (21%)	94 (94%)	5 (5%)	1 (1%)	13	23
8	T	51/835 (6%)	51 (100%)	0	0	100	100
All	All	910/2303 (40%)	897 (99%)	12 (1%)	1 (0%)	50	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	O	237	LYS

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	85/111 (77%)	84 (99%)	1 (1%)	67	84
1	E	85/111 (77%)	85 (100%)	0	100	100
2	B	68/79 (86%)	68 (100%)	0	100	100
2	F	68/79 (86%)	67 (98%)	1 (2%)	60	80
3	C	85/100 (85%)	85 (100%)	0	100	100
3	G	92/100 (92%)	92 (100%)	0	100	100
4	D	81/105 (77%)	81 (100%)	0	100	100
4	H	81/105 (77%)	81 (100%)	0	100	100
7	O	94/369 (26%)	92 (98%)	2 (2%)	48	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	T	45/633 (7%)	45 (100%)	0	100	100
All	All	784/1792 (44%)	780 (100%)	4 (0%)	85	94

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	LYS
2	F	22	LEU
7	O	203	LEU
7	O	233	ARG

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

Mol	Chain	Res	Type
1	E	85	GLN
8	T	293	ASN
8	T	302	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

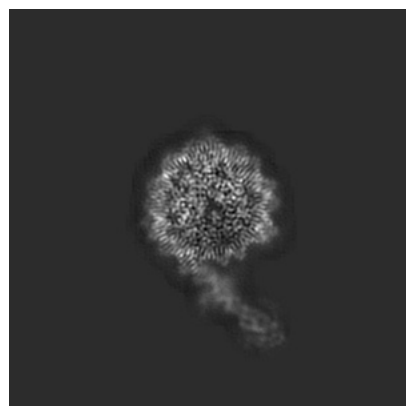
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-43197. 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](#)

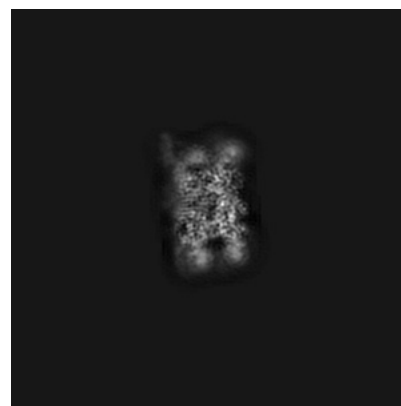
6.1.1 Primary map



X

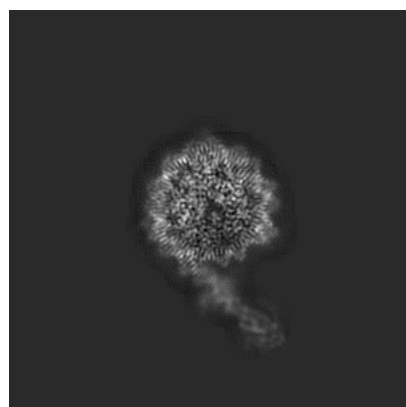


Y



Z

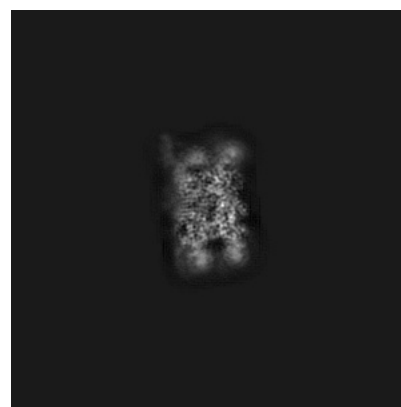
6.1.2 Raw map



X



Y

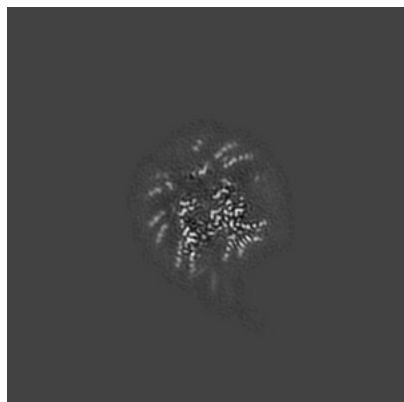


Z

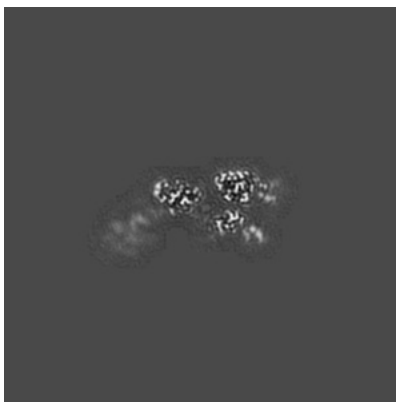
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

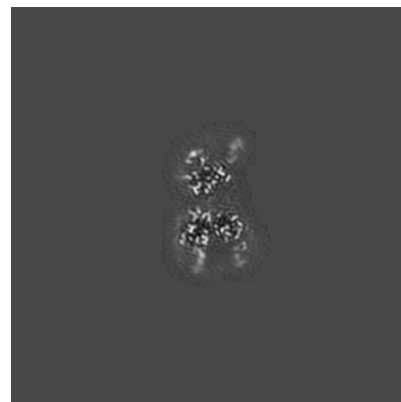
6.2.1 Primary map



X Index: 150

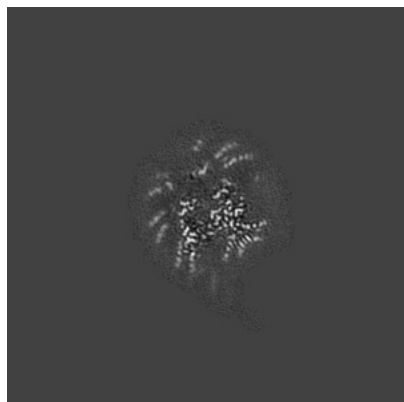


Y Index: 150

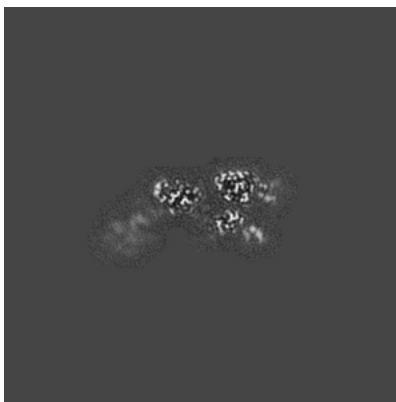


Z Index: 150

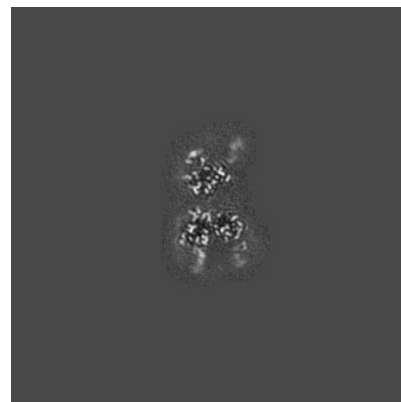
6.2.2 Raw map



X Index: 150



Y Index: 150

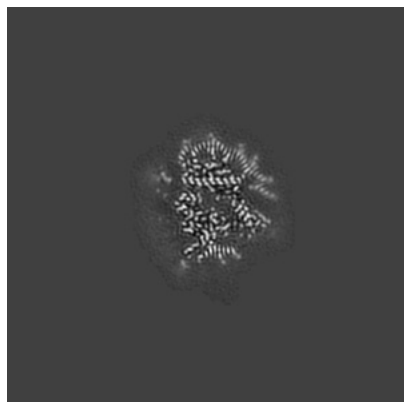


Z Index: 150

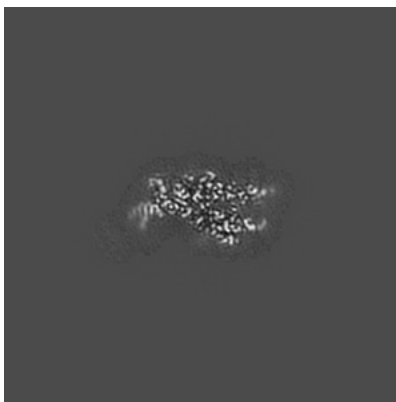
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

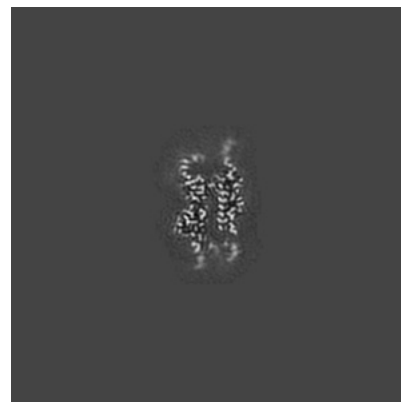
6.3.1 Primary map



X Index: 158

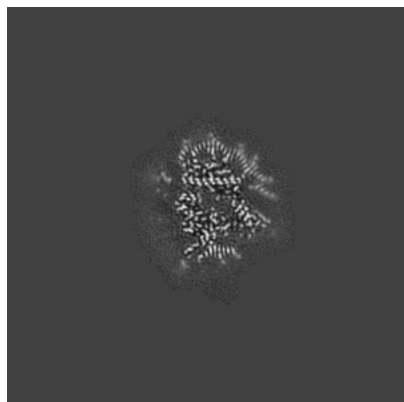


Y Index: 138

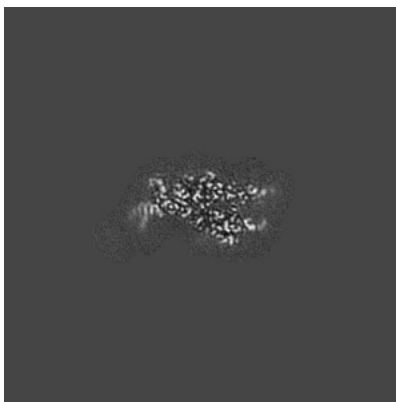


Z Index: 167

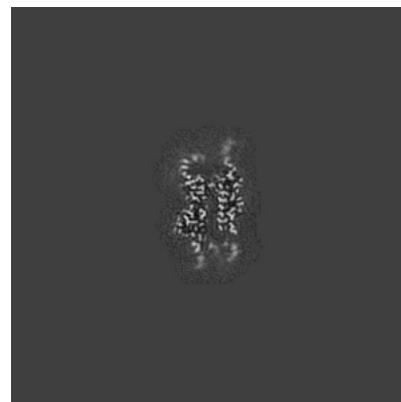
6.3.2 Raw map



X Index: 158



Y Index: 138

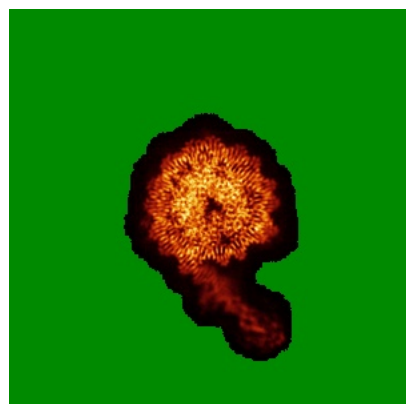


Z Index: 167

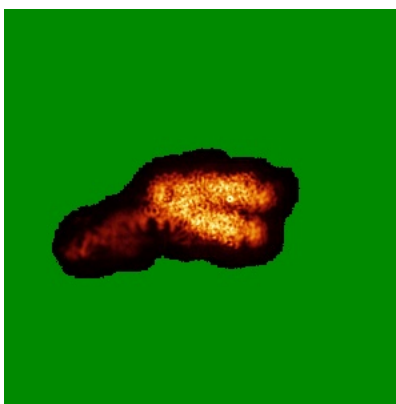
The images above show the largest variance slices of the map in three orthogonal directions.

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

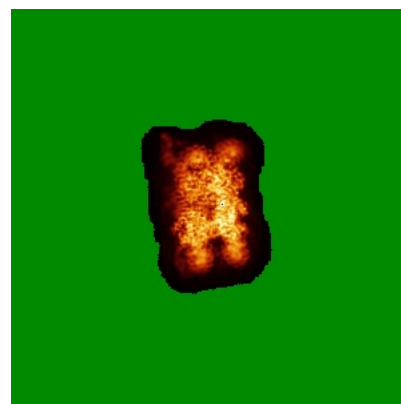
6.4.1 Primary map



X

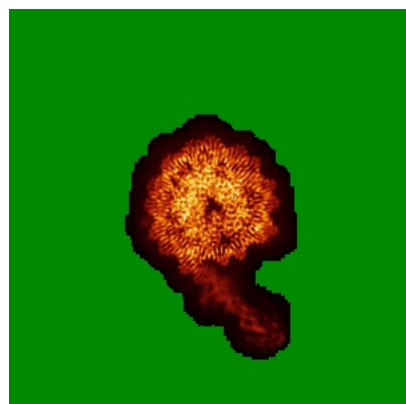


Y

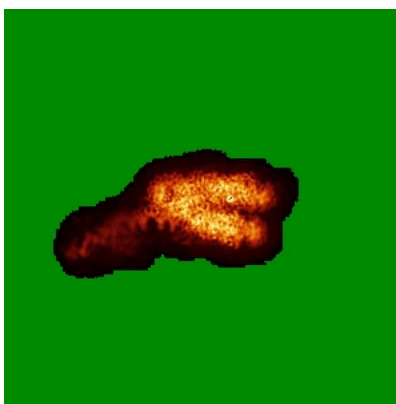


Z

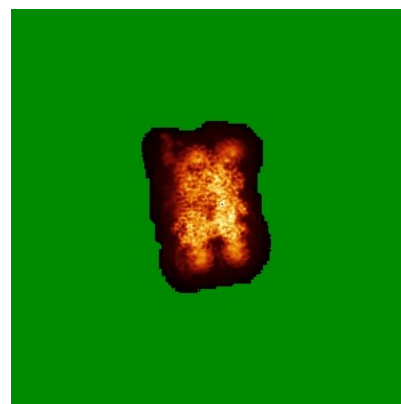
6.4.2 Raw map



X



Y



Z

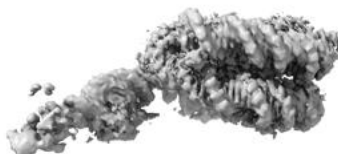
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

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 0.004. 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.

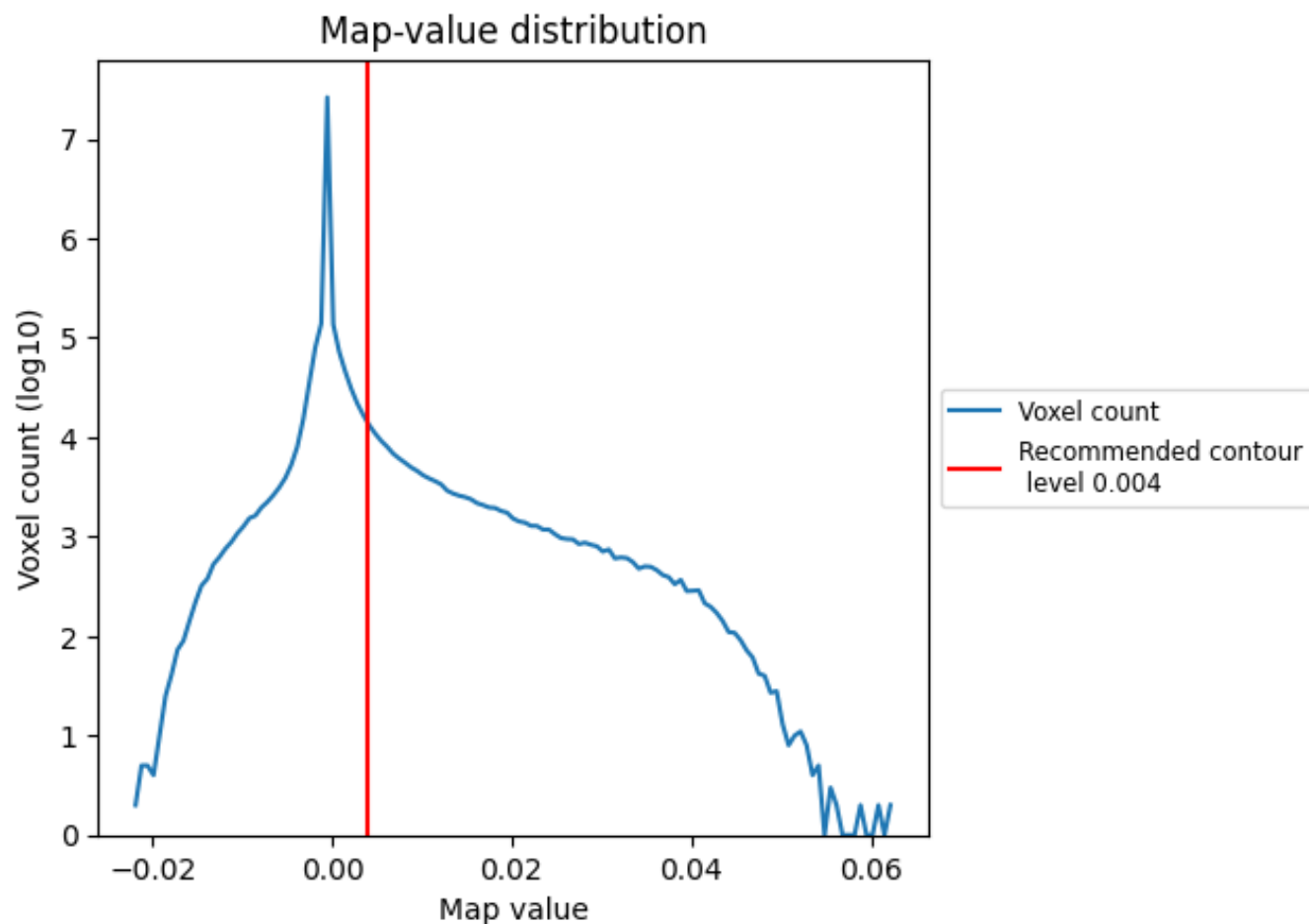
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

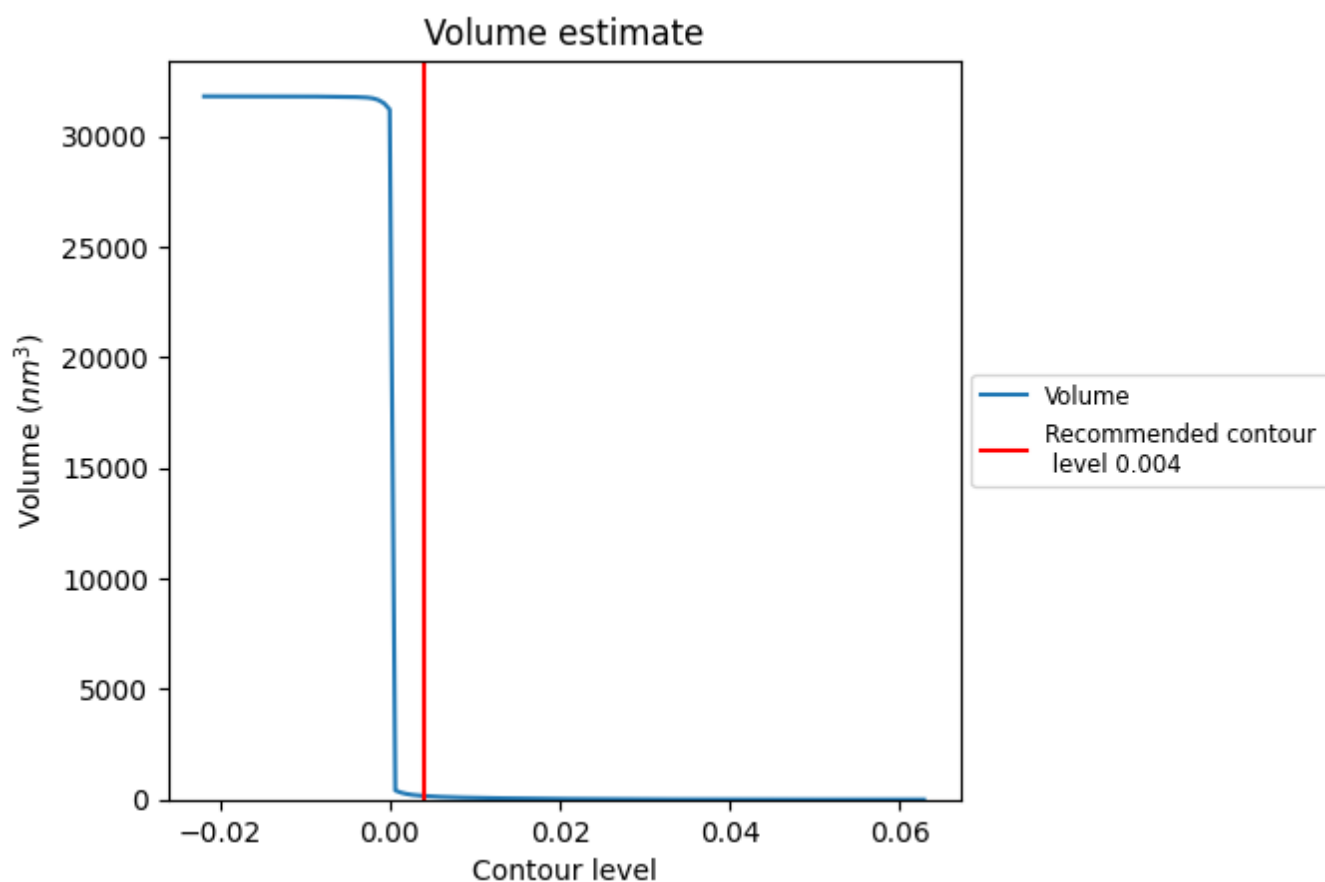
This section contains the results of statistical analysis of the map.

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.

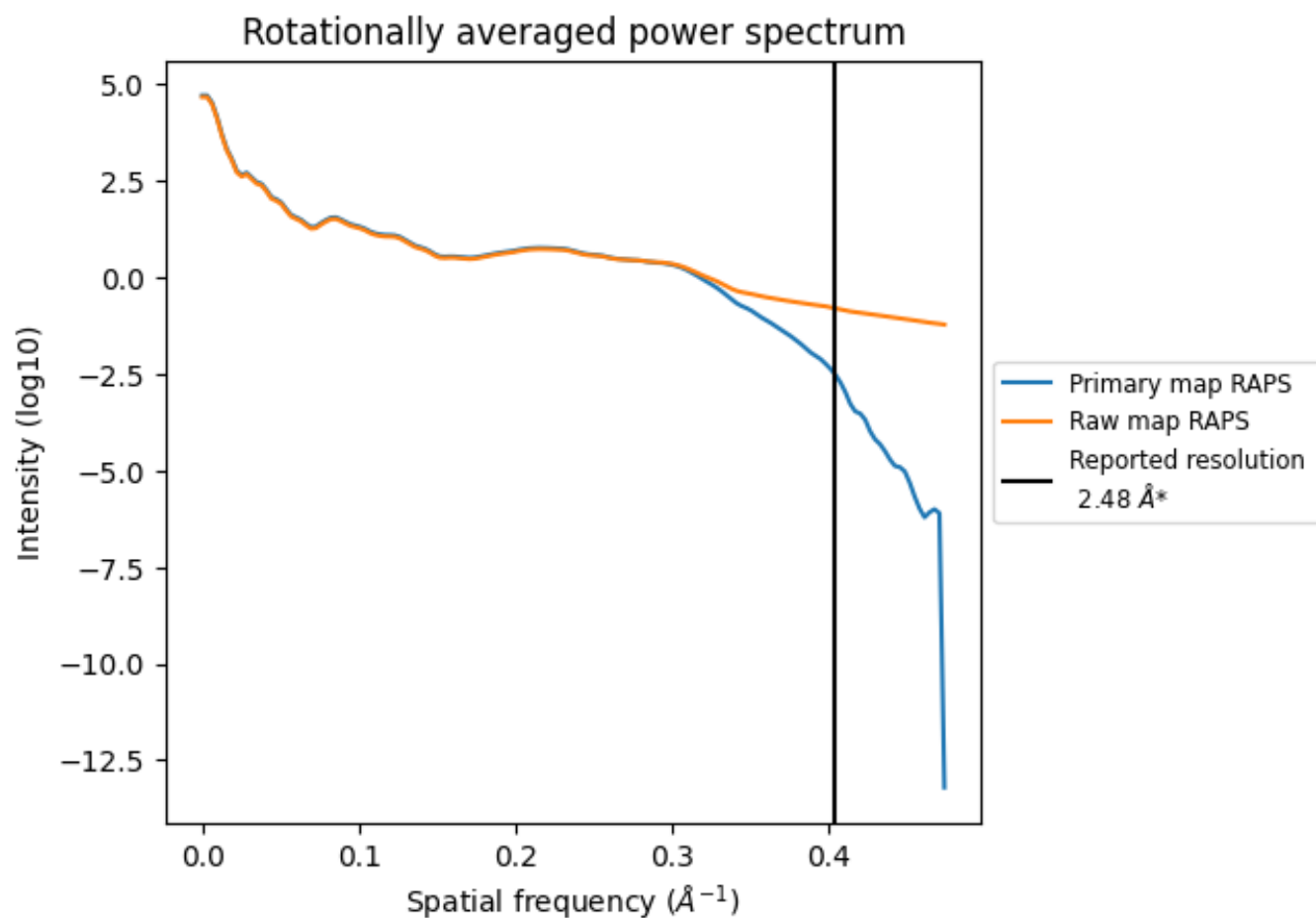
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 163 nm^3 ; this corresponds to an approximate mass of 148 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 ⓘ

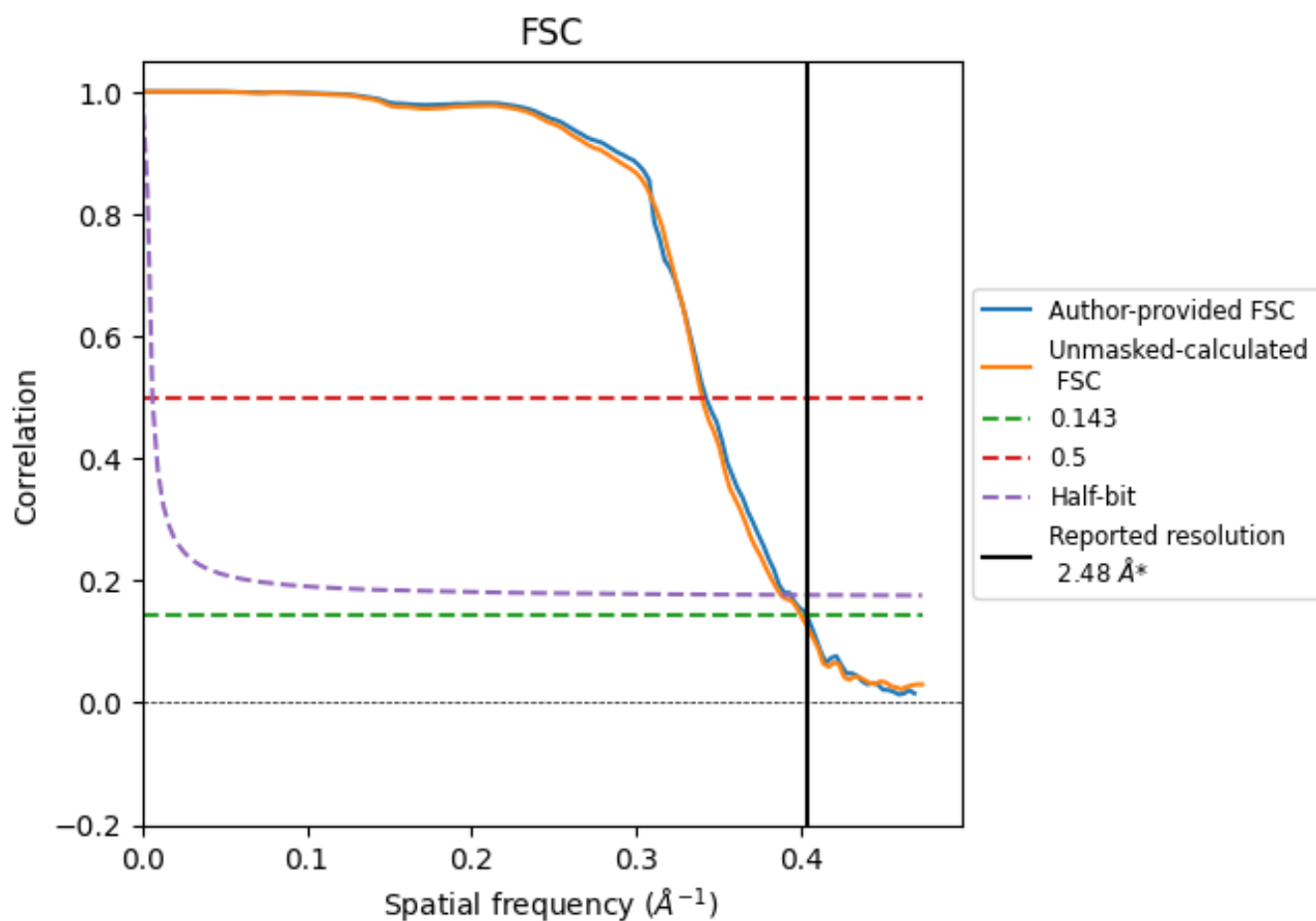


*Reported resolution corresponds to spatial frequency of 0.403 Å⁻¹

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 0.403 \AA^{-1}

8.2 Resolution estimates [i](#)

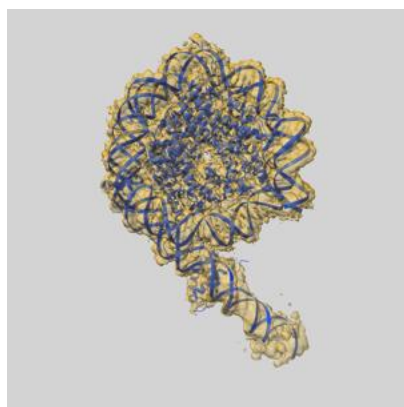
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.48	-	-
Author-provided FSC curve	2.48	2.92	2.54
Unmasked-calculated*	2.50	2.94	2.58

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

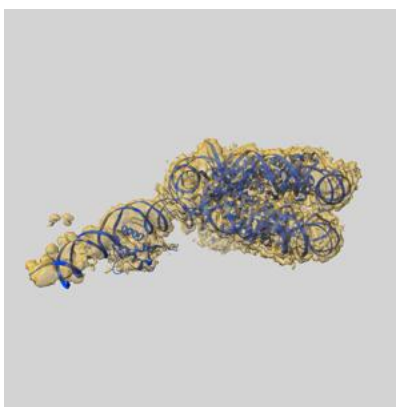
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-43197 and PDB model 8VG1. Per-residue inclusion information can be found in section [3](#) on page [7](#).

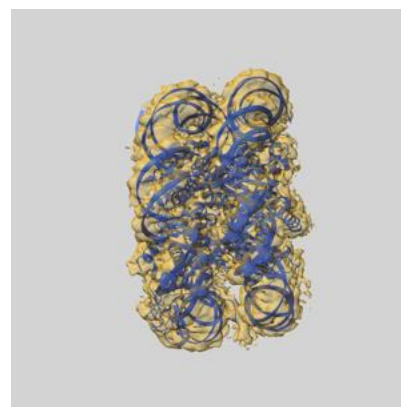
9.1 Map-model overlay [i](#)



X



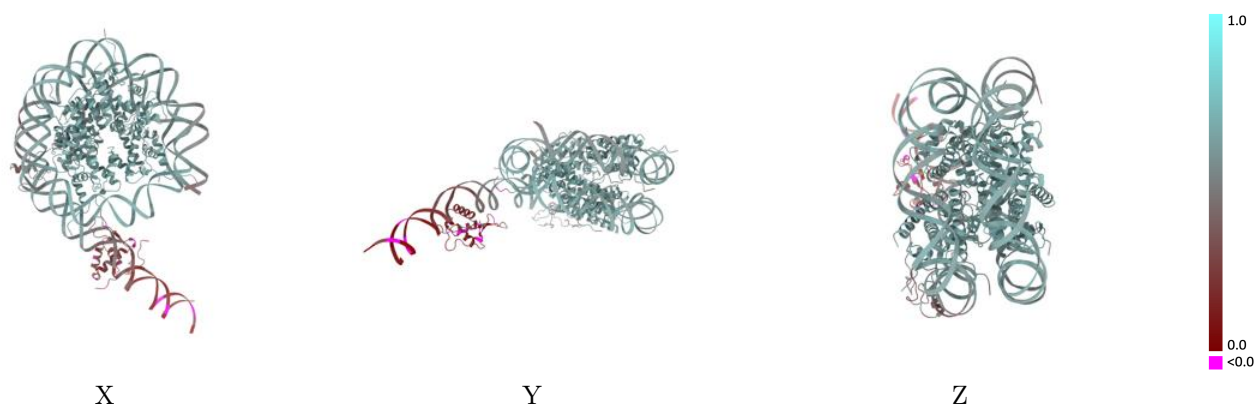
Y



Z

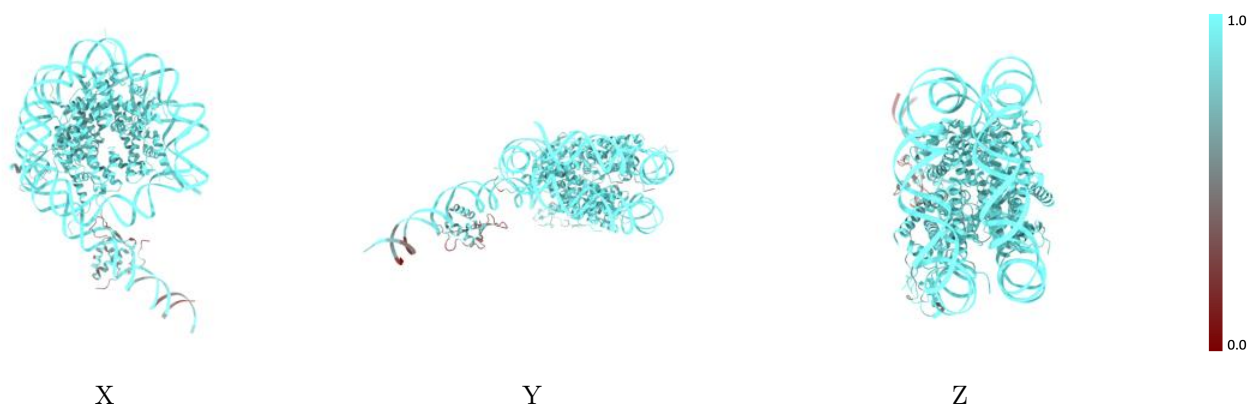
The images above show the 3D surface view of the map at the recommended contour level 0.004 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



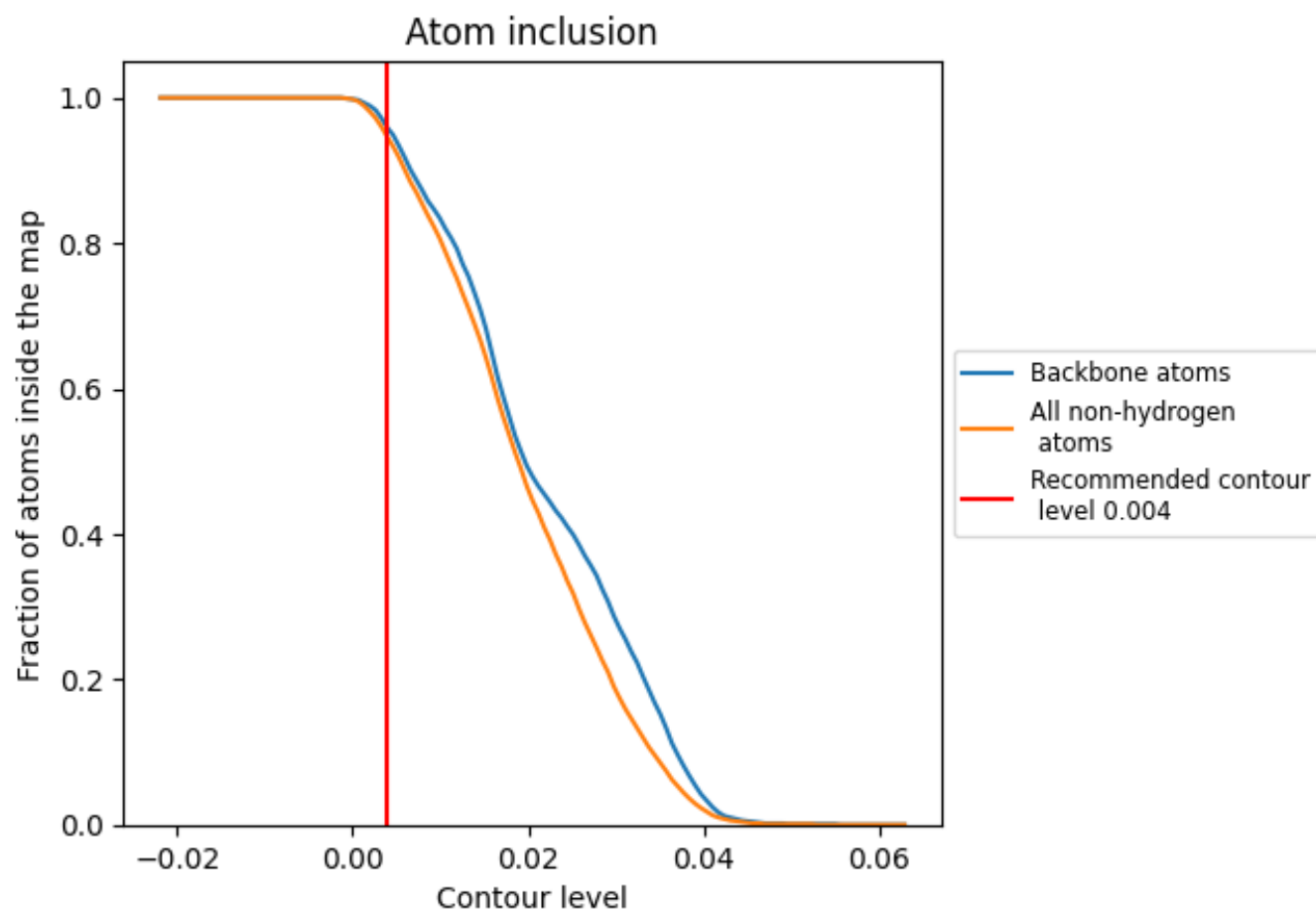
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.004).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 94% 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 (0.004) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.9450	<div></div> 0.5410
A	<div></div> 0.9870	<div></div> 0.6340
B	<div></div> 0.9870	<div></div> 0.6360
C	<div></div> 0.9870	<div></div> 0.6240
D	<div></div> 0.9730	<div></div> 0.6100
E	<div></div> 0.9920	<div></div> 0.6370
F	<div></div> 0.9830	<div></div> 0.6290
G	<div></div> 0.9530	<div></div> 0.6000
H	<div></div> 0.9850	<div></div> 0.6170
I	<div></div> 0.9670	<div></div> 0.5230
J	<div></div> 0.9630	<div></div> 0.5240
O	<div></div> 0.5910	<div></div> 0.1520
T	<div></div> 0.8160	<div></div> 0.4270

1.0

0.0

<0.0