



wwPDB EM Validation Summary Report ⓘ

Mar 7, 2026 – 11:54 PM UTC

PDB ID : 8SJD / pdb_00008sjd
EMDB ID : EMD-40553
Title : Cryo-EM structure of the Hermes transposase bound to two right-ends of its DNA transposon.
Authors : Lannes, L.; Dyda, F.
Deposited on : 2023-04-17
Resolution : 5.10 Å (reported)
Based on initial models : 4D1Q, 8EB5, 6DX0

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.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

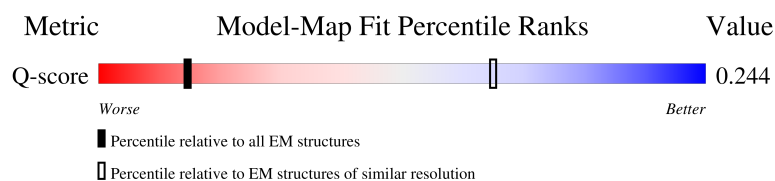
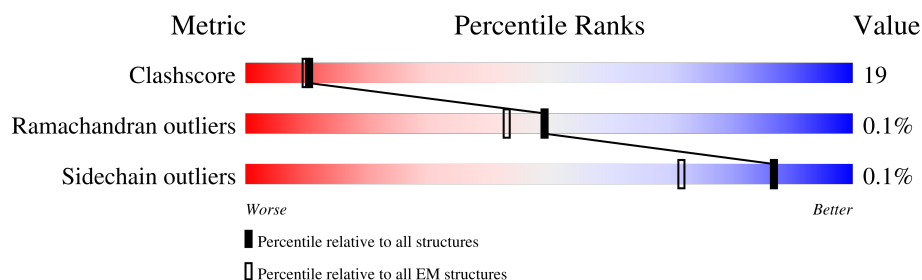
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 5.10 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.



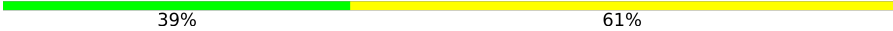
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	844 (4.60 - 5.60)

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	612	10% 44% 34% 22%
1	B	612	10% 46% 32% 22%
1	C	612	27% 42% 35% 23%
1	D	612	34% 48% 30% 22%

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Mol	Chain	Length	Quality of chain
2	G	55	
2	J	55	
3	F	46	
3	I	46	
4	E	7	
4	H	7	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19777 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 Hermes transposase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	474	Total	C	N	O	S	0	0
			3817	2443	645	710	19		
1	D	479	Total	C	N	O	S	0	0
			3857	2467	653	718	19		
1	B	477	Total	C	N	O	S	0	0
			3838	2459	648	712	19		
1	A	478	Total	C	N	O	S	0	0
			3846	2463	650	714	19		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	2	GLU	GLN	engineered mutation	UNP Q25438
C	128	GLY	LYS	engineered mutation	UNP Q25438
D	2	GLU	GLN	engineered mutation	UNP Q25438
D	128	GLY	LYS	engineered mutation	UNP Q25438
B	2	GLU	GLN	engineered mutation	UNP Q25438
B	128	GLY	LYS	engineered mutation	UNP Q25438
A	2	GLU	GLN	engineered mutation	UNP Q25438
A	128	GLY	LYS	engineered mutation	UNP Q25438

- Molecule 2 is a DNA chain called DNA (55-MER).

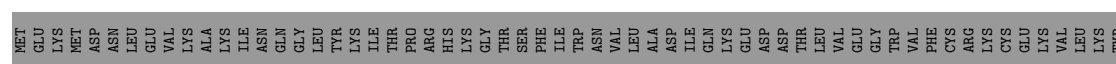
Mol	Chain	Residues	Atoms					AltConf	Trace
2	J	55	Total	C	N	O	P	0	0
			1119	538	182	345	54		
2	G	55	Total	C	N	O	P	0	0
			1122	538	182	347	55		

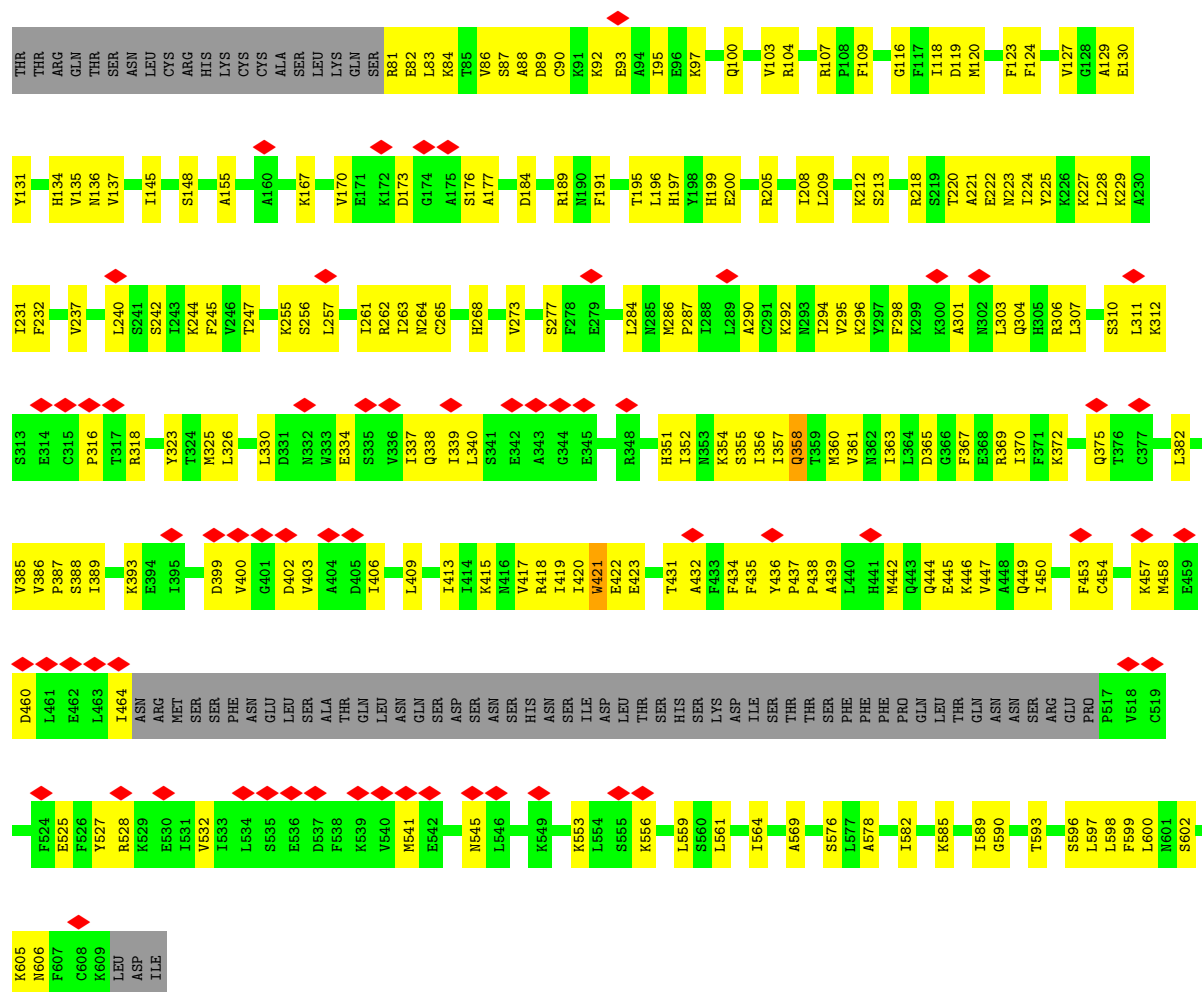
- Molecule 3 is a DNA chain called DNA (46-MER).

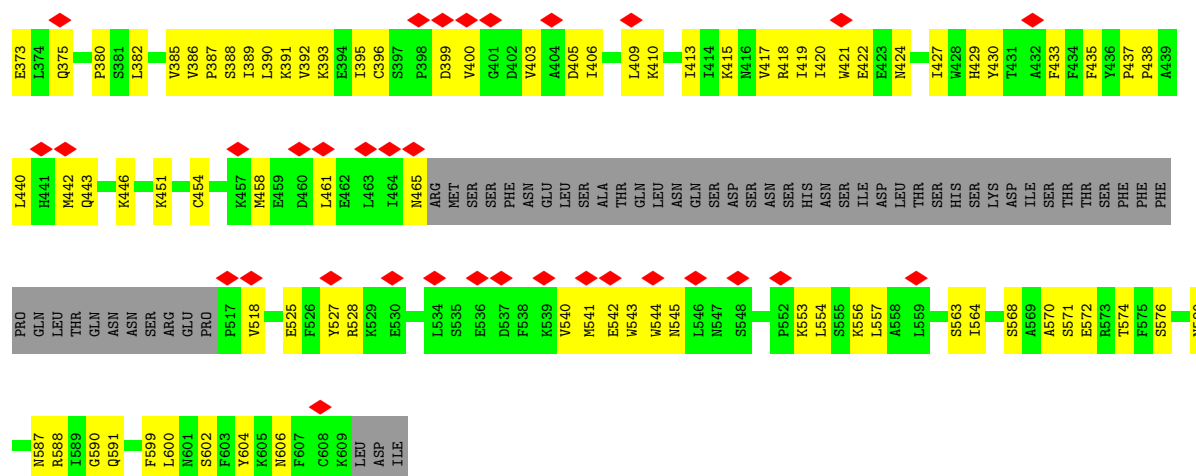
Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	46	Total	C	N	O	P	0	0
			945	449	193	258	45		
3	F	46	Total	C	N	O	P	0	0
			945	449	193	258	45		

- Molecule 4 is a DNA chain called DNA (8-MER).

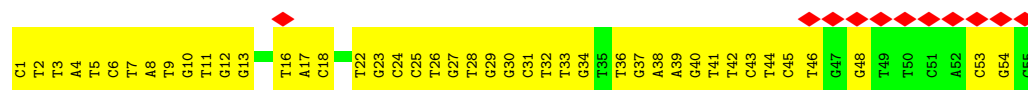
Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	7	Total	C	N	O	P	0	0
			144	69	30	39	6		
4	E	7	Total	C	N	O	P	0	0
			144	69	30	39	6		



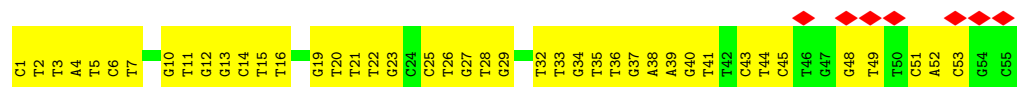




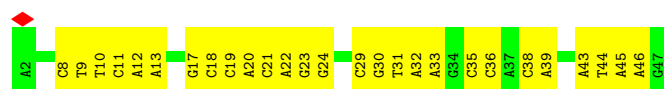
- Molecule 2: DNA (55-MER)



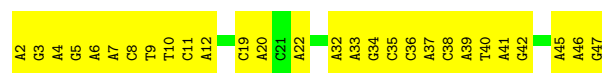
- Molecule 2: DNA (55-MER)



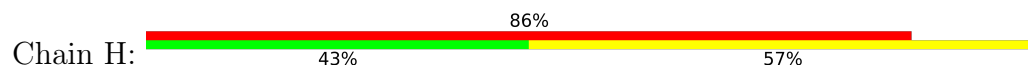
- Molecule 3: DNA (46-MER)

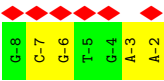


- Molecule 3: DNA (46-MER)

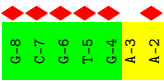
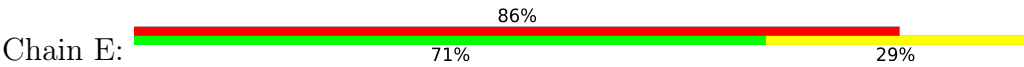


- Molecule 4: DNA (8-MER)





● Molecule 4: DNA (8-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	53656	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	48.7	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.045	Depositor
Minimum map value	-0.020	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	249.40001, 249.40001, 249.40001	wwPDB
Map dimensions	290, 290, 290	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.86, 0.86, 0.86	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	0.17	0/3919	0.38	0/5287
1	B	0.18	0/3911	0.42	2/5276 (0.0%)
1	C	0.14	0/3889	0.35	0/5246
1	D	0.14	0/3931	0.34	0/5304
2	G	0.22	0/1251	0.44	0/1930
2	J	0.23	0/1248	0.46	0/1926
3	F	0.22	0/1066	0.33	0/1642
3	I	0.22	0/1066	0.36	0/1642
4	E	0.19	0/162	0.31	0/249
4	H	0.18	0/162	0.39	0/249
All	All	0.17	0/20605	0.38	2/28751 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	421	TRP	CA-C-N	-6.50	109.32	122.31
1	B	421	TRP	C-N-CA	-6.50	109.32	122.31

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	3846	0	3916	168	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3838	0	3910	156	0
1	C	3817	0	3884	171	0
1	D	3857	0	3922	141	0
2	G	1122	0	629	44	0
2	J	1119	0	630	50	0
3	F	945	0	513	28	0
3	I	945	0	513	22	0
4	E	144	0	80	2	0
4	H	144	0	80	2	0
All	All	19777	0	18077	710	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 19.

The worst 5 of 710 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:C:118:ILE:HG22	1:C:122:LYS:HZ3	1.30	0.94
1:B:367:PHE:HA	1:B:370:ILE:HD12	1.61	0.83
2:J:6:DC:H2''	2:J:7:DT:H5'	1.66	0.78
1:B:205:ARG:HH21	1:B:606:ASN:HB3	1.48	0.78
1:A:465:ASN:HD21	1:A:556:LYS:HE2	1.48	0.78

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	474/612 (78%)	444 (94%)	30 (6%)	0	100	100
1	B	473/612 (77%)	442 (93%)	31 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	470/612 (77%)	441 (94%)	29 (6%)	0	100	100
1	D	475/612 (78%)	449 (94%)	25 (5%)	1 (0%)	43	78
All	All	1892/2448 (77%)	1776 (94%)	115 (6%)	1 (0%)	49	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	518	VAL

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	437/565 (77%)	437 (100%)	0	100	100
1	B	436/565 (77%)	435 (100%)	1 (0%)	87	86
1	C	434/565 (77%)	434 (100%)	0	100	100
1	D	439/565 (78%)	439 (100%)	0	100	100
All	All	1746/2260 (77%)	1745 (100%)	1 (0%)	87	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	358	GLN

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

Mol	Chain	Res	Type
1	A	293	ASN
1	A	465	ASN
1	A	580	ASN
1	B	260	ASN
1	B	362	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 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.

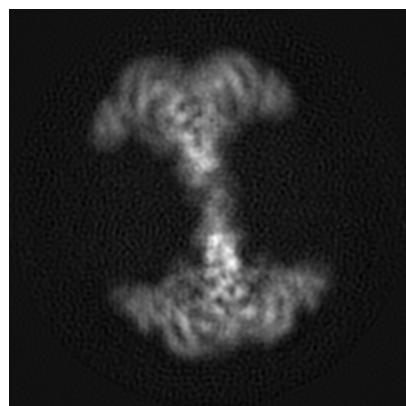
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40553. These allow visual inspection of the internal detail of the map and identification of artifacts.

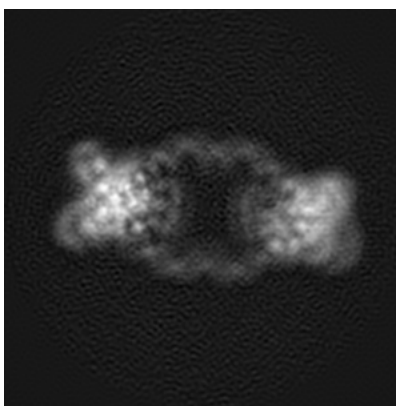
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

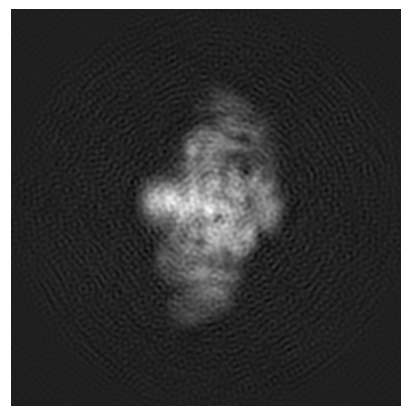
6.1.1 Primary map



X

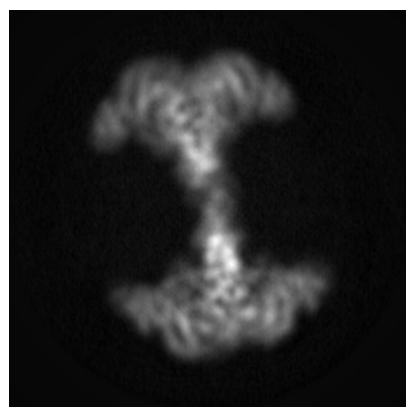


Y

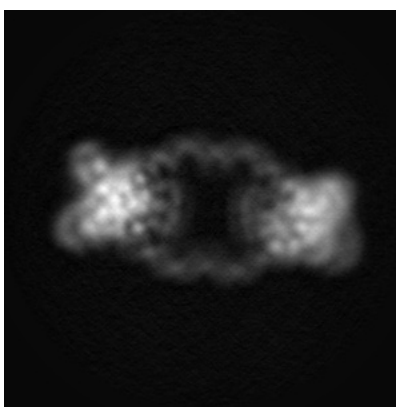


Z

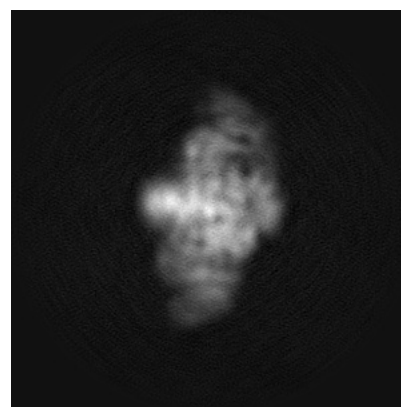
6.1.2 Raw map



X



Y

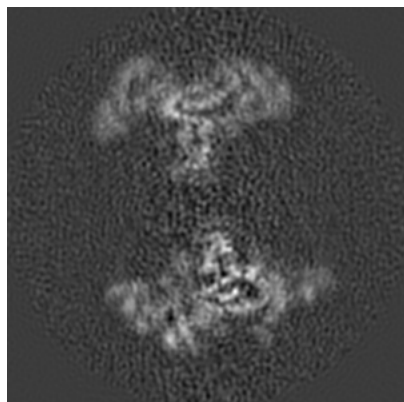


Z

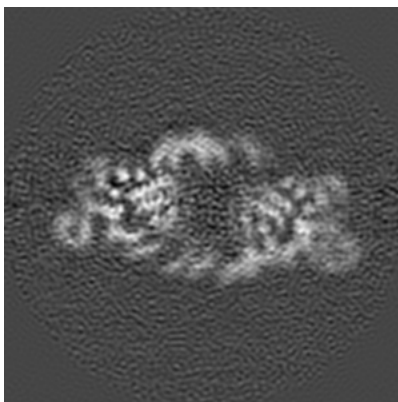
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

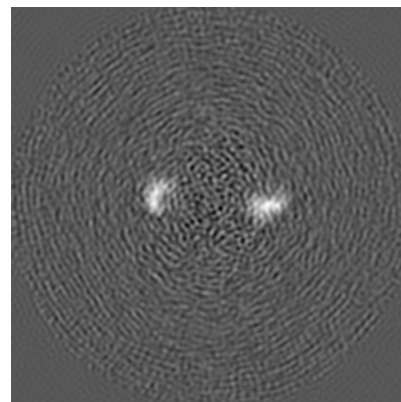
6.2.1 Primary map



X Index: 145

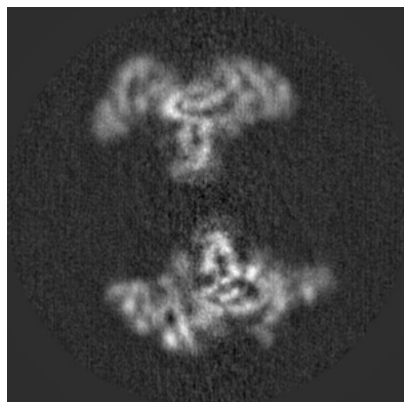


Y Index: 145

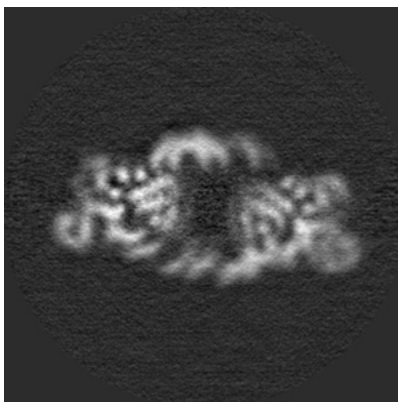


Z Index: 145

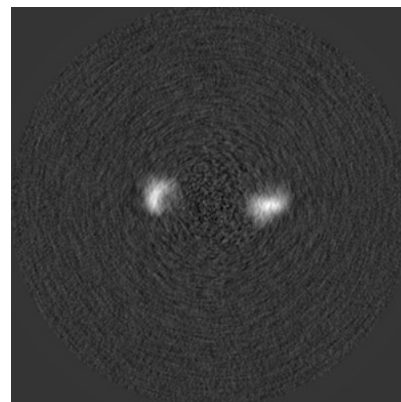
6.2.2 Raw map



X Index: 145



Y Index: 145

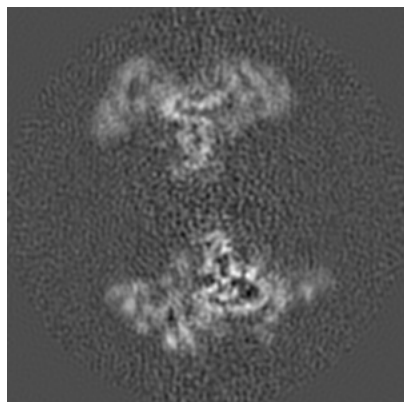


Z Index: 145

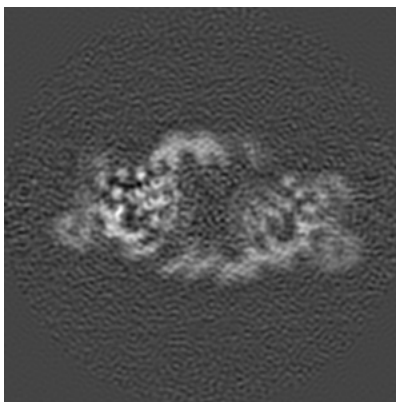
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

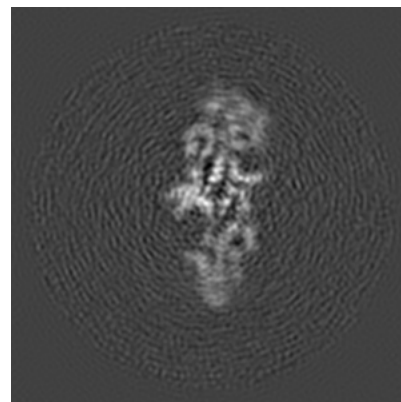
6.3.1 Primary map



X Index: 144

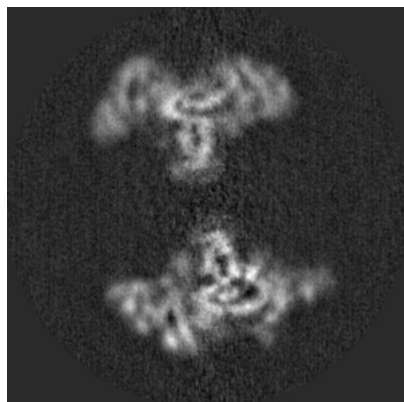


Y Index: 147

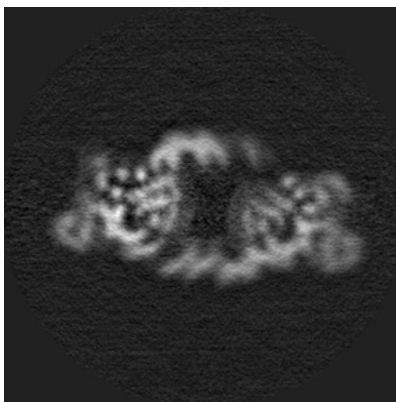


Z Index: 83

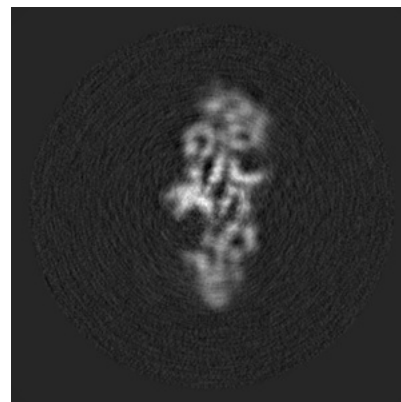
6.3.2 Raw map



X Index: 144



Y Index: 147

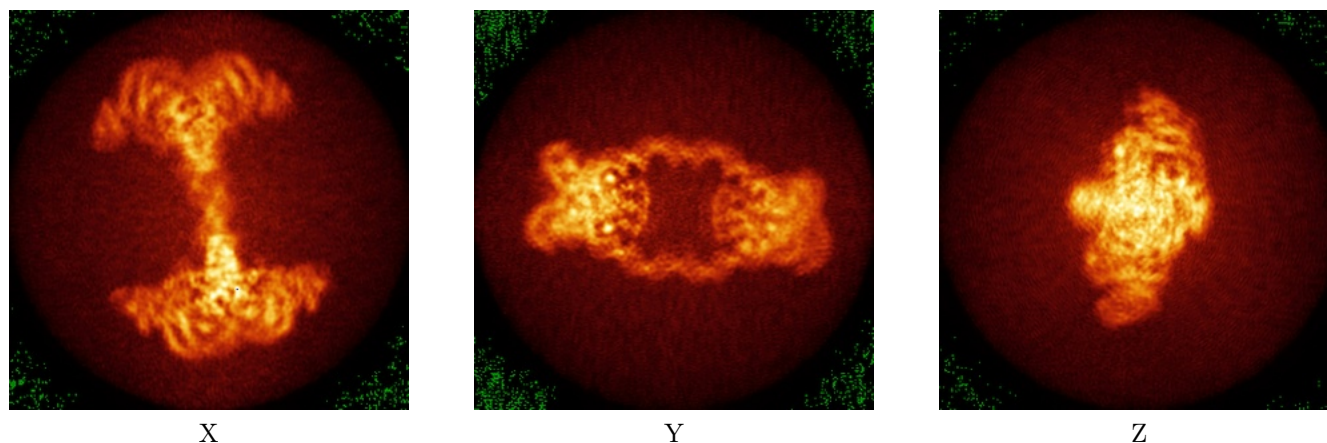


Z Index: 83

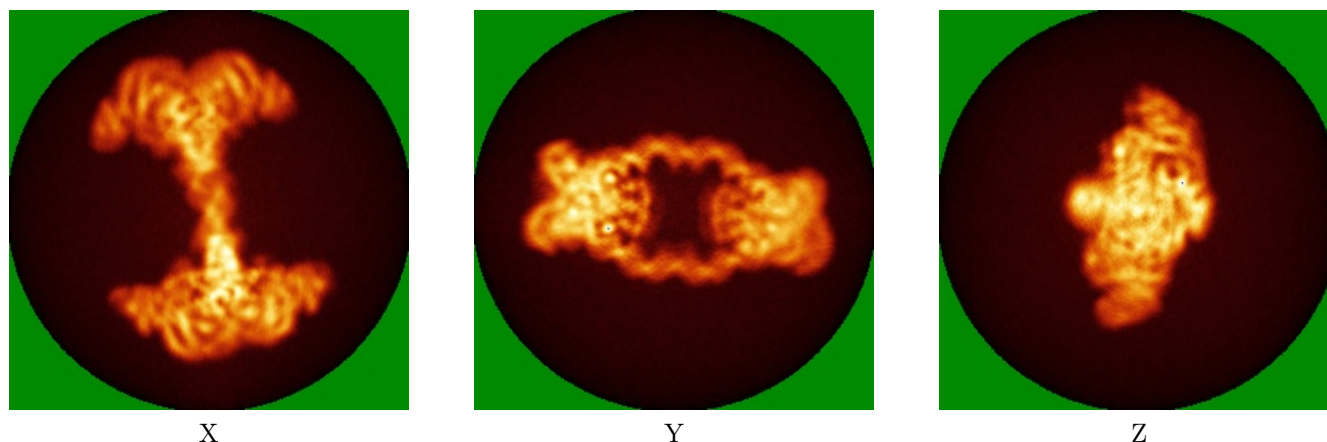
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

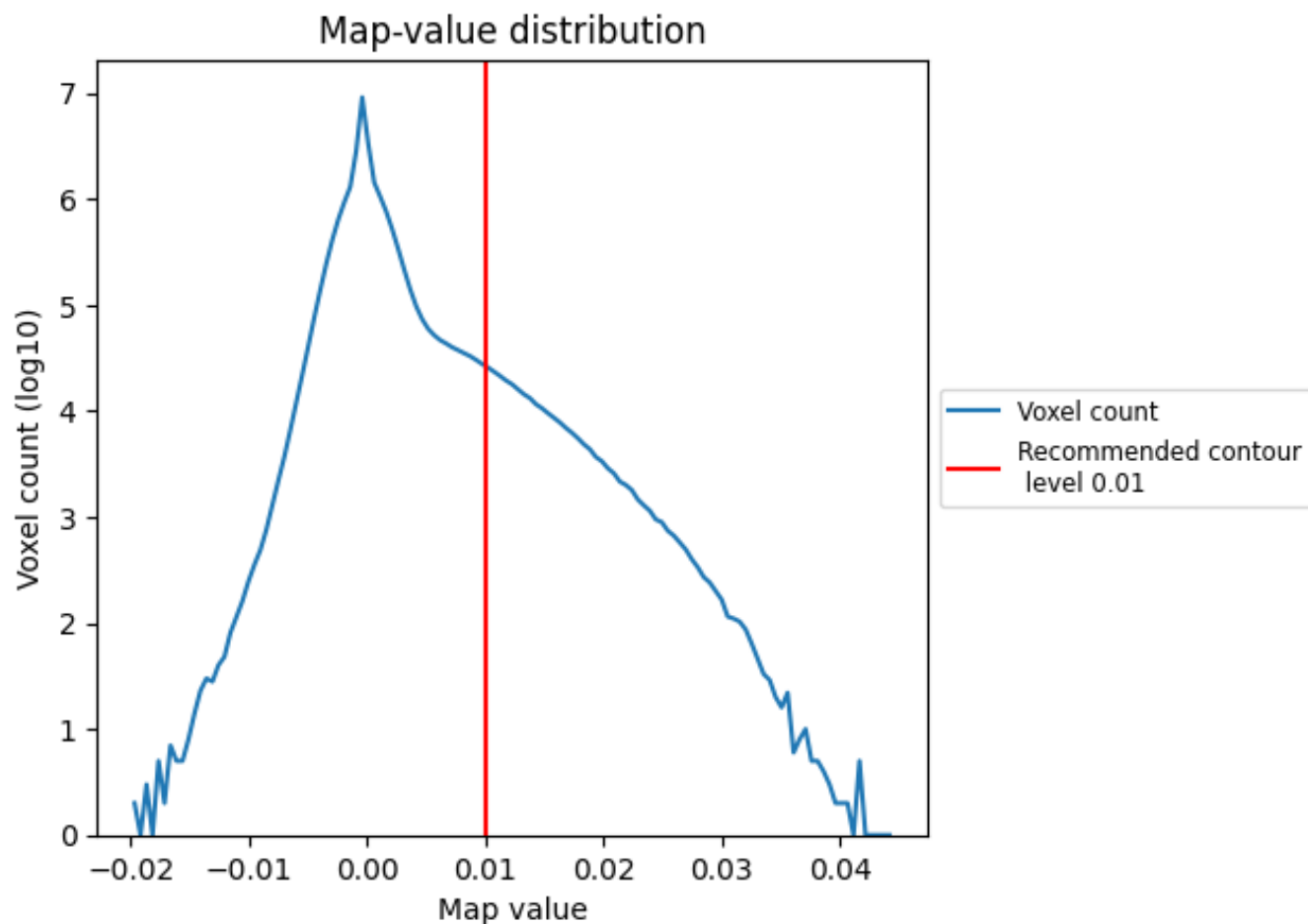
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

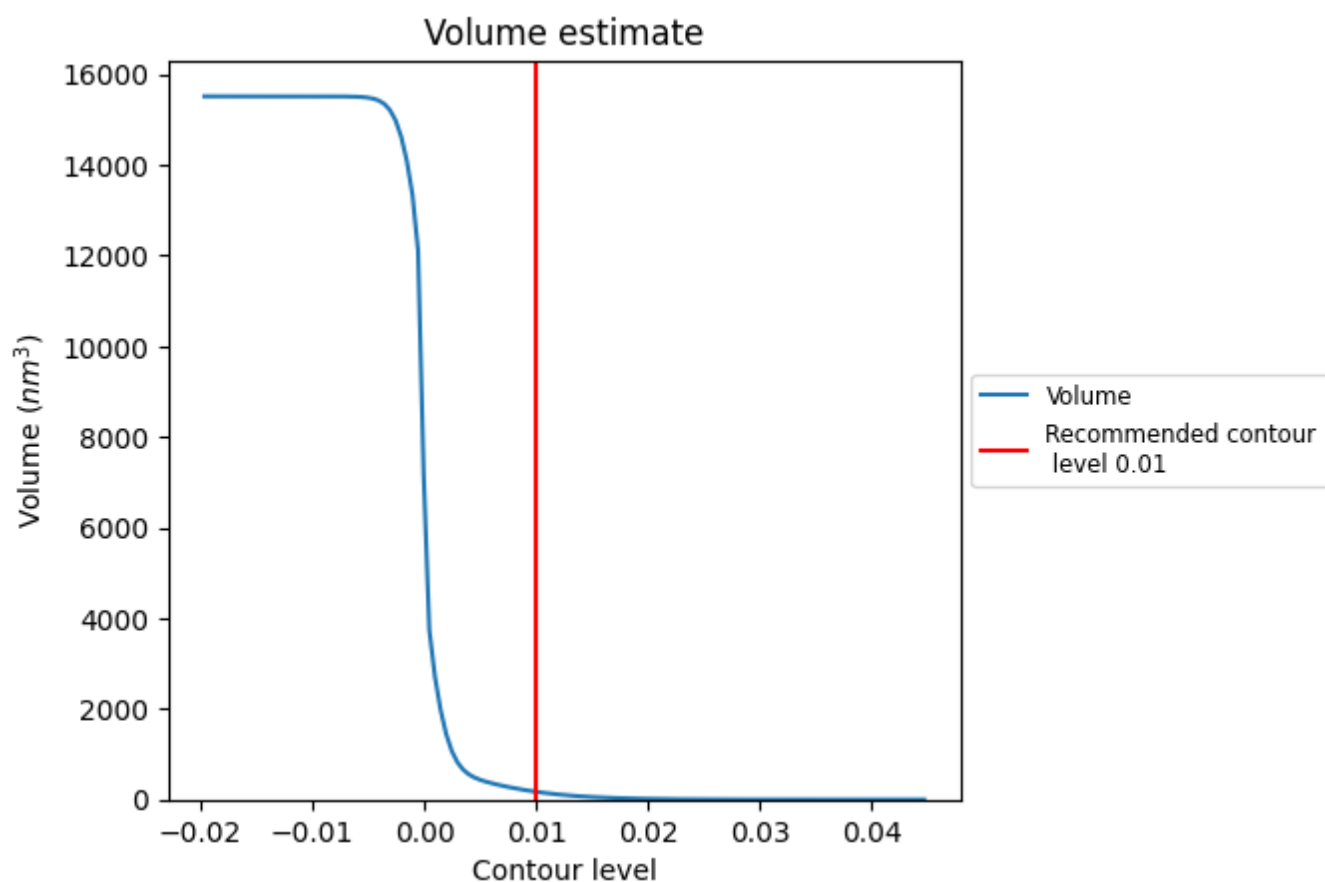
This section contains the results of statistical analysis of the map.

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.

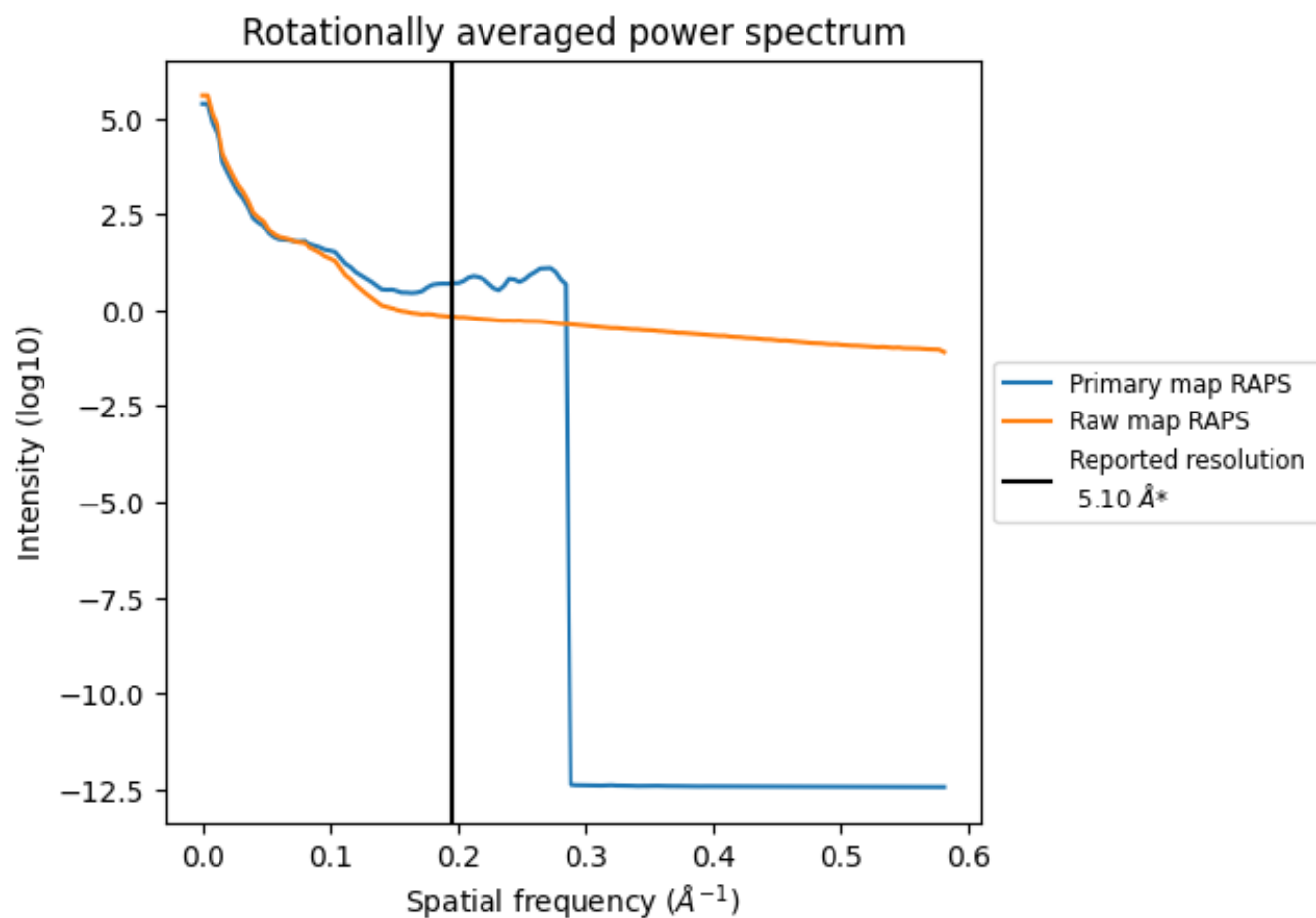
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 172 nm³; this corresponds to an approximate mass of 155 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 ⓘ

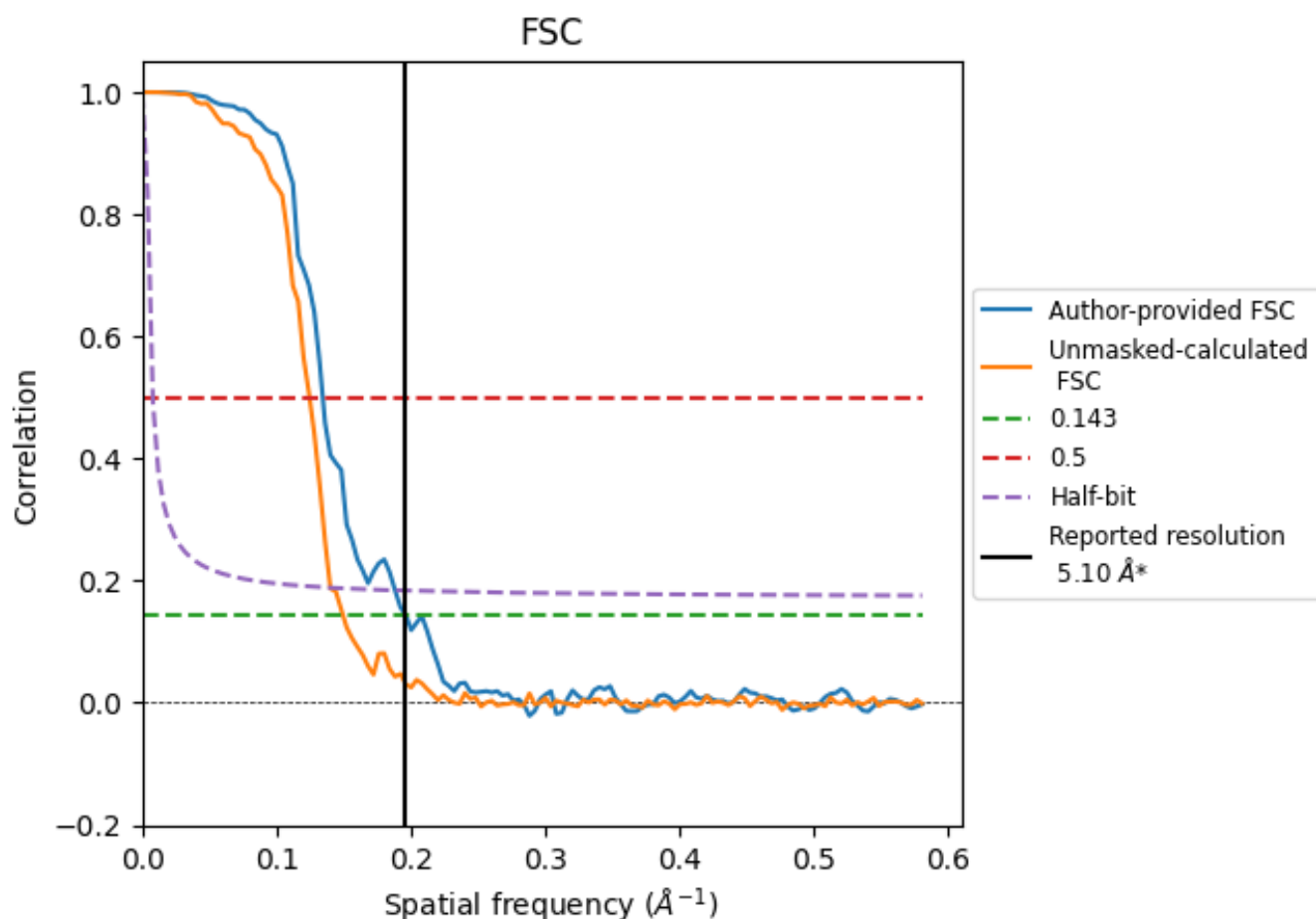


*Reported resolution corresponds to spatial frequency of 0.196 Å⁻¹

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 $0.196 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	5.10	-	-
Author-provided FSC curve	5.09	7.43	5.31
Unmasked-calculated*	6.68	8.02	7.13

*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 6.68 differs from the reported value 5.1 by more than 10 %

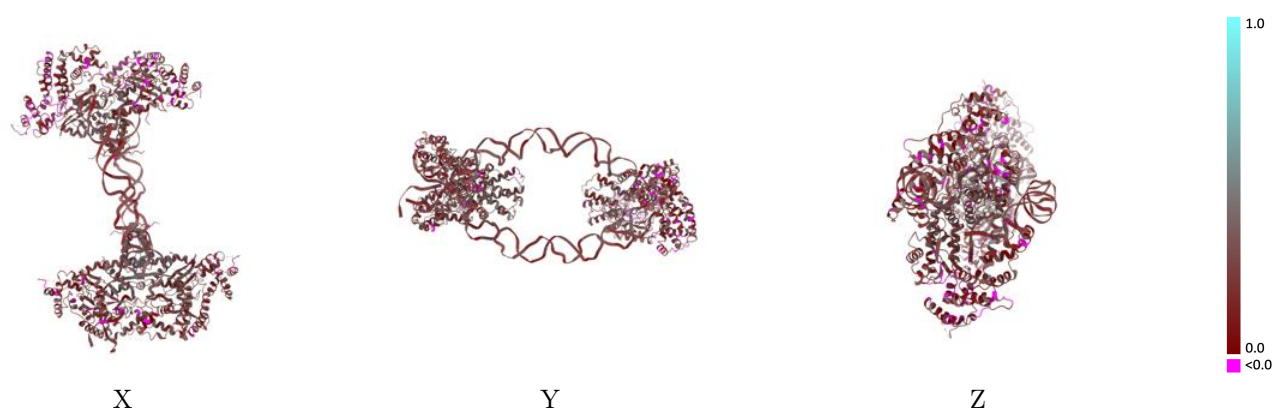
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-40553 and PDB model 8SJD. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

This section was not generated.

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

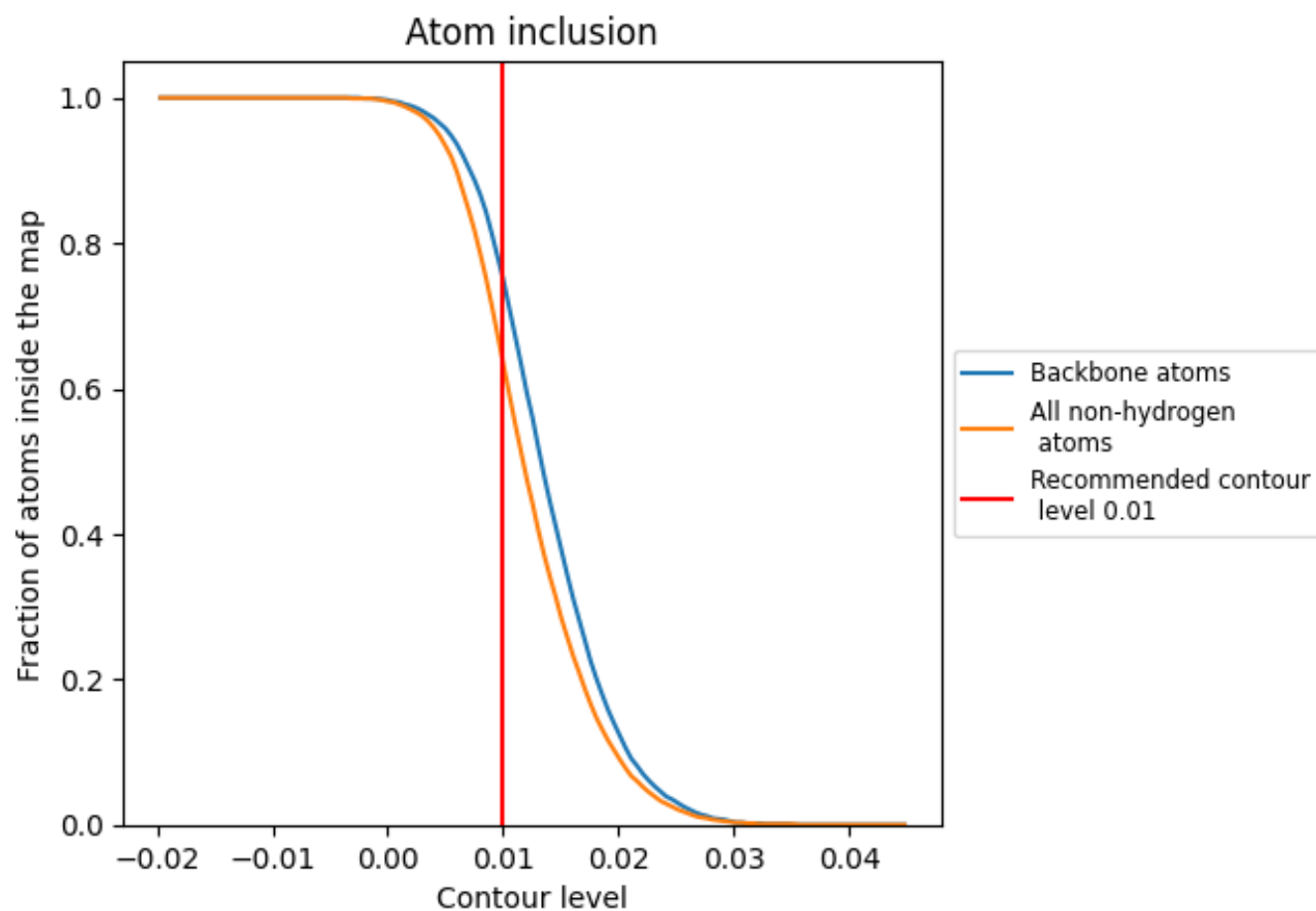


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 64% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6350	0.2440
A	0.6970	0.2680
B	0.6960	0.2700
C	0.5200	0.2260
D	0.4770	0.1980
E	0.1390	0.2850
F	0.8760	0.2710
G	0.7740	0.2600
H	0.1530	0.2270
I	0.8550	0.2540
J	0.7340	0.2460

1.0

0.0

<0.0