



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 06:34 PM UTC

PDB ID : 9IK9 / pdb_00009ik9
EMDB ID : EMD-60651
Title : Cryo-EM Structure of SST analogs bound SSTR1-Gi complex
Authors : Wong, T.S.; Zeng, Z.C.; Xiong, T.T.; Gan, S.Y.; Du, Y.
Deposited on : 2024-06-26
Resolution : 3.37 Å (reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

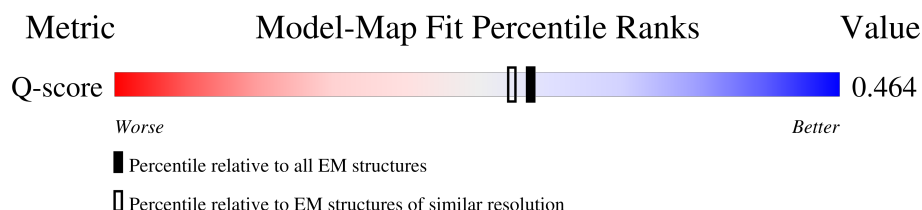
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : **FAILED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

i

ELECTRON MICROSCOPY

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	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Q-score	25397	14287 (2.87 - 3.87)

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	4J2	F	1	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17030 atoms, of which 8425 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 Guanine nucleotide-binding protein G(i) subunit alpha-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	217	Total	C	H	N	O	S	0	0
			3261	1065	1597	286	301	12		

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	336	Total	C	H	N	O	S	0	0
			4915	1564	2391	450	491	19		

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-21	MET	-	initiating methionine	UNP P62873
B	-20	HIS	-	expression tag	UNP P62873
B	-19	HIS	-	expression tag	UNP P62873
B	-18	HIS	-	expression tag	UNP P62873
B	-17	HIS	-	expression tag	UNP P62873
B	-16	HIS	-	expression tag	UNP P62873
B	-15	HIS	-	expression tag	UNP P62873
B	-14	HIS	-	expression tag	UNP P62873
B	-13	HIS	-	expression tag	UNP P62873
B	-12	HIS	-	expression tag	UNP P62873
B	-11	HIS	-	expression tag	UNP P62873
B	-10	LEU	-	expression tag	UNP P62873
B	-9	GLU	-	expression tag	UNP P62873
B	-8	VAL	-	expression tag	UNP P62873
B	-7	LEU	-	expression tag	UNP P62873
B	-6	PHE	-	expression tag	UNP P62873
B	-5	GLN	-	expression tag	UNP P62873
B	-4	GLY	-	expression tag	UNP P62873
B	-3	PRO	-	expression tag	UNP P62873
B	-2	GLY	-	expression tag	UNP P62873

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	SER	-	expression tag	UNP P62873
B	0	SER	-	expression tag	UNP P62873
B	1	GLY	-	expression tag	UNP P62873
B	341	VAL	-	expression tag	UNP P62873
B	342	SER	-	expression tag	UNP P62873
B	343	GLY	-	expression tag	UNP P62873
B	344	TRP	-	expression tag	UNP P62873
B	345	ARG	-	expression tag	UNP P62873
B	346	LEU	-	expression tag	UNP P62873
B	347	PHE	-	expression tag	UNP P62873
B	348	LYS	-	expression tag	UNP P62873
B	349	LYS	-	expression tag	UNP P62873
B	350	ILE	-	expression tag	UNP P62873
B	351	SER	-	expression tag	UNP P62873

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	53	Total	C	H	N	O	S	0	0
			749	241	372	66	67	3		

- Molecule 4 is a protein called scFv16.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	E	232	Total	C	H	N	O	S	0	0
			3416	1116	1668	295	328	9		

- Molecule 5 is a protein called Somatostatin receptor type 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	D	283	Total	C	H	N	O	S	0	0
			4543	1467	2328	361	363	24		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	392	LEU	-	expression tag	UNP P30872
D	393	GLU	-	expression tag	UNP P30872

- Molecule 6 is a protein (with D amino acids) called (4J2)(DCY)(DTY)(DTR)K(DVA)(DCY)(ALO)(NH2).

Mol	Chain	Residues	Atoms						AltConf	Trace
6	F	9	Total	C	H	N	O	S	0	1
			146	54	69	11	10	2		

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

3 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	217201	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 ($e^-/\text{\AA}^2$)	53.6	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	1400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.833	Depositor
Minimum map value	-1.171	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	326.40002, 326.40002, 326.40002	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8500001, 0.8500001, 0.8500001	Depositor

4 Model quality [i](#)

4.1 Standard geometry [i](#)

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

4.2 Too-close contacts [i](#)

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

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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

4.3.2 Protein sidechains [i](#)

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

4.3.3 RNA [i](#)

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

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

7 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	ALO	F	8	6	5,6,7	0.71	0	5,7,9	1.01	1 (20%)
6	DTY	F	3	6	11,12,13	4.75	6 (54%)	10,15,17	0.81	0
6	DVA	F	6	6	4,6,7	0.51	0	6,7,9	0.56	0
6	DCY	F	2	6	4,5,6	0.58	0	1,5,7	1.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	DCY	F	7	6	4,5,6	0.92	0	1,5,7	1.73	0
6	4J2	F	1	6	15,16,17	8.57	10 (66%)	16,21,23	5.63	15 (93%)
6	DTR	F	4	6	14,15,16	2.72	5 (35%)	15,20,22	4.82	8 (53%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	ALO	F	8	6	-	4/5/6/8	-
6	DTY	F	3	6	-	2/5/6/8	0/1/1/1
6	DVA	F	6	6	-	0/5/6/8	-
6	DCY	F	2	6	-	1/1/4/6	-
6	DCY	F	7	6	-	1/1/4/6	-
6	4J2	F	1	6	-	2/5/6/8	0/2/2/2
6	DTR	F	4	6	-	0/5/6/8	0/2/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	1	4J2	CD1-CG	-16.83	1.05	1.38
6	F	1	4J2	CZ3-CZ2	13.13	1.64	1.36
6	F	1	4J2	CE3-CE2	-13.11	1.13	1.42
6	F	1	4J2	CE1-CD1	-10.92	1.13	1.36
6	F	3	DTY	CE1-CD1	9.44	1.54	1.38
6	F	1	4J2	CE4-CE3	-9.26	1.17	1.36
6	F	1	4J2	CZ2-CZ1	8.70	1.61	1.42
6	F	1	4J2	CD2-CE2	8.63	1.61	1.42
6	F	1	4J2	CE2-CZ1	8.39	1.61	1.42
6	F	3	DTY	CE2-CZ	7.25	1.52	1.39
6	F	3	DTY	CD2-CG	7.00	1.52	1.38
6	F	1	4J2	CD2-CG	-6.15	1.23	1.37
6	F	4	DTR	CD2-CE2	5.55	1.48	1.41
6	F	4	DTR	CD2-CG	5.18	1.52	1.44
6	F	4	DTR	CD1-NE1	-4.59	1.29	1.37
6	F	3	DTY	CE2-CD2	-4.20	1.32	1.38
6	F	3	DTY	CD1-CG	-4.18	1.30	1.38
6	F	3	DTY	CE1-CZ	-3.97	1.31	1.39
6	F	4	DTR	CB-CG	3.01	1.55	1.50
6	F	4	DTR	CD1-CG	-2.38	1.33	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	1	4J2	CB-CG	2.05	1.56	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	4	DTR	CE2-NE1-CD1	13.30	120.91	109.08
6	F	1	4J2	CZ2-CZ1-CE1	10.17	146.18	123.01
6	F	1	4J2	CD2-CE2-CZ1	-7.81	108.13	118.96
6	F	1	4J2	CZ2-CZ1-CE2	-7.58	105.89	118.91
6	F	4	DTR	CD2-CE2-NE1	-7.46	101.31	107.52
6	F	1	4J2	CE1-CZ1-CE2	-6.39	107.94	118.91
6	F	1	4J2	CD1-CG-CD2	5.92	129.66	118.88
6	F	1	4J2	CD1-CE1-CZ1	5.92	129.68	120.84
6	F	1	4J2	CE4-CE3-CE2	5.77	128.50	120.48
6	F	1	4J2	CZ3-CE4-CE3	5.70	128.03	120.40
6	F	1	4J2	CE3-CE2-CZ1	5.58	128.48	118.91
6	F	4	DTR	CG-CD1-NE1	-5.22	104.58	110.31
6	F	4	DTR	CE2-CD2-CG	-4.90	102.21	107.17
6	F	4	DTR	CD2-CG-CD1	4.72	110.99	106.16
6	F	1	4J2	CE4-CZ3-CZ2	-4.47	114.41	120.40
6	F	1	4J2	CZ3-CZ2-CZ1	-4.17	114.68	120.48
6	F	4	DTR	CZ2-CE2-NE1	4.00	137.14	130.74
6	F	1	4J2	CE1-CD1-CG	3.91	128.71	121.07
6	F	4	DTR	CB-CG-CD2	-3.53	119.92	126.70
6	F	4	DTR	CE3-CD2-CG	3.46	139.02	133.85
6	F	1	4J2	CB-CG-CD2	-3.03	114.20	120.91
6	F	1	4J2	CG-CD2-CE2	-2.86	115.89	121.28
6	F	1	4J2	CB-CG-CD1	-2.56	116.14	120.90
6	F	8	ALO	O-C-CA	-2.12	119.31	124.77

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	1	4J2	N-CA-CB-CG
6	F	1	4J2	C-CA-CB-CG
6	F	3	DTY	C-CA-CB-CG
6	F	7	DCY	N-CA-CB-SG
6	F	8	ALO	N-CA-CB-CG2
6	F	8	ALO	C-CA-CB-CG2
6	F	8	ALO	C-CA-CB-OG1
6	F	2	DCY	N-CA-CB-SG

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Mol	Chain	Res	Type	Atoms
6	F	3	DTY	N-CA-CB-CG
6	F	8	ALO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

4.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

4.6 Ligand geometry [i](#)

There are no ligands in this entry.

4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

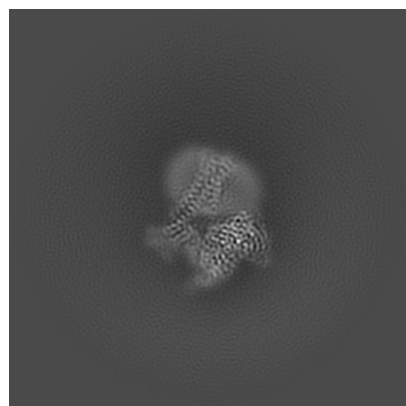
5 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-60651. These allow visual inspection of the internal detail of the map and identification of artifacts.

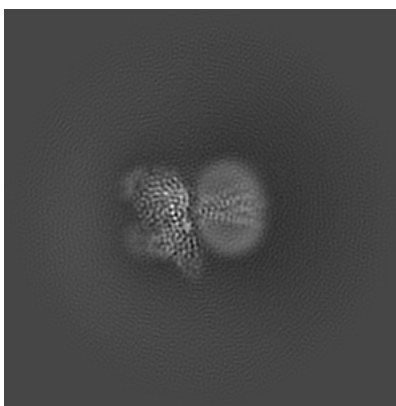
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

5.1 Orthogonal projections [i](#)

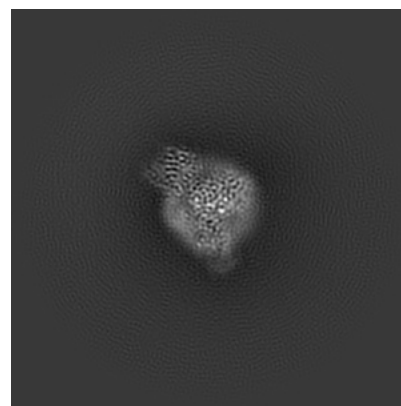
5.1.1 Primary map



X

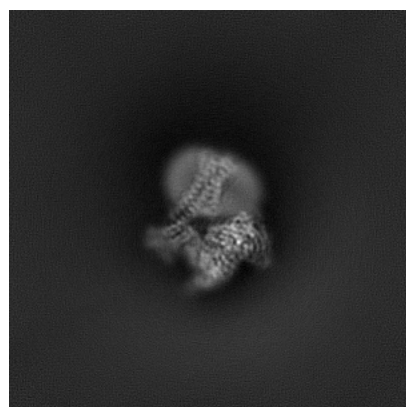


Y

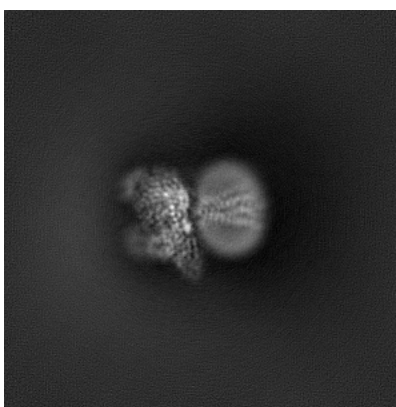


Z

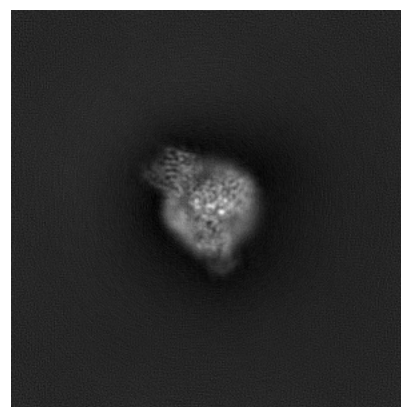
5.1.2 Raw map



X



Y

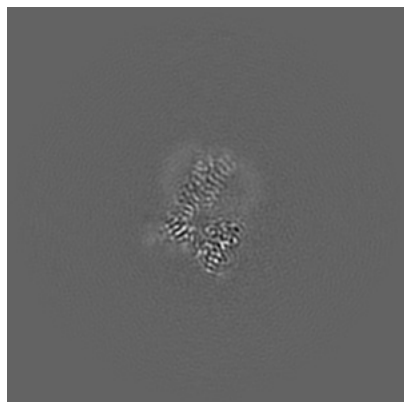


Z

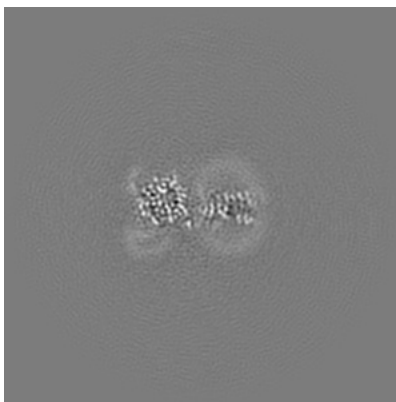
The images above show the map projected in three orthogonal directions.

5.2 Central slices [i](#)

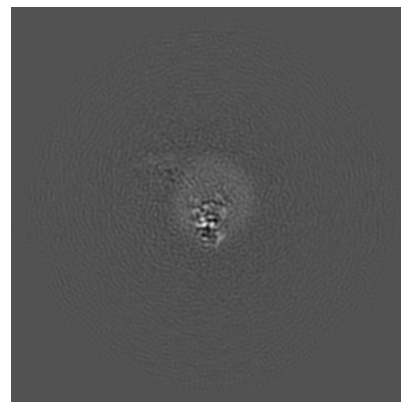
5.2.1 Primary map



X Index: 192

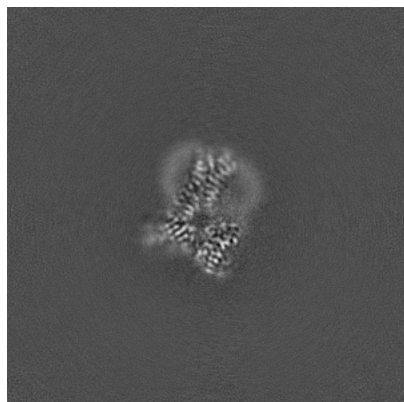


Y Index: 192

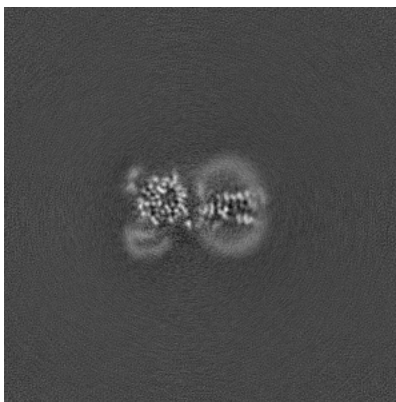


Z Index: 192

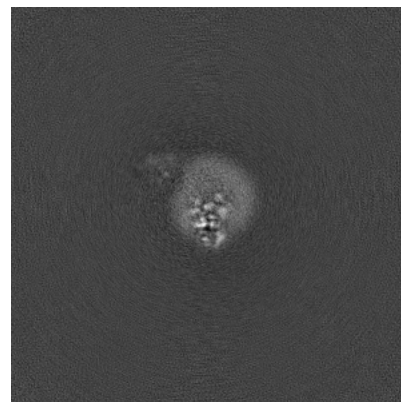
5.2.2 Raw map



X Index: 192



Y Index: 192

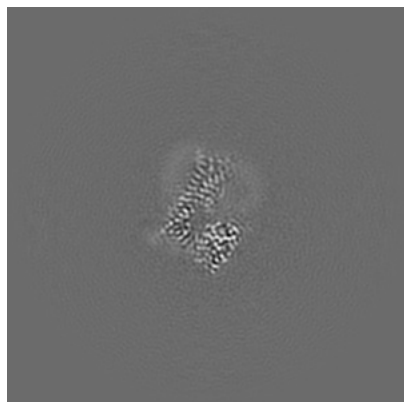


Z Index: 192

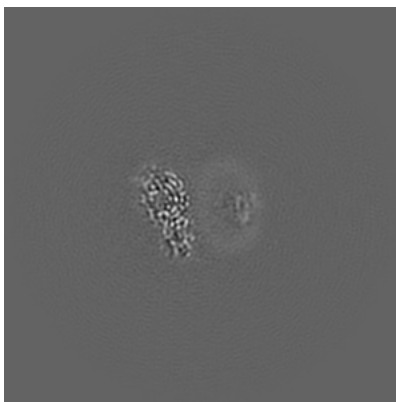
The images above show central slices of the map in three orthogonal directions.

5.3 Largest variance slices [i](#)

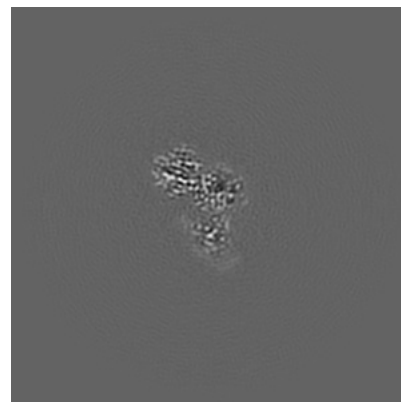
5.3.1 Primary map



X Index: 188

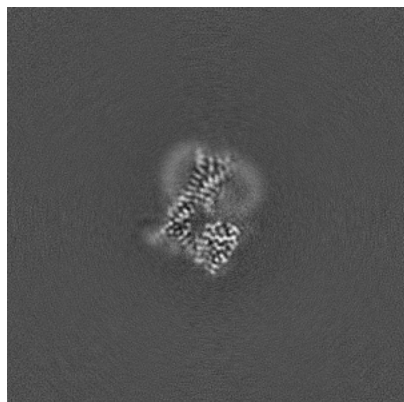


Y Index: 209

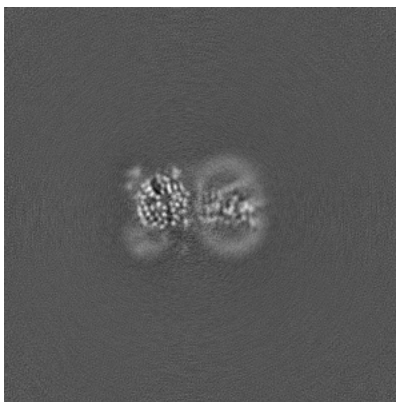


Z Index: 167

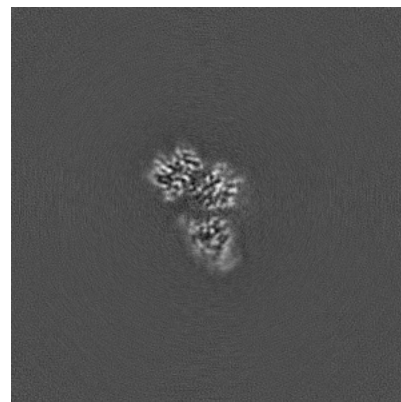
5.3.2 Raw map



X Index: 188



Y Index: 196

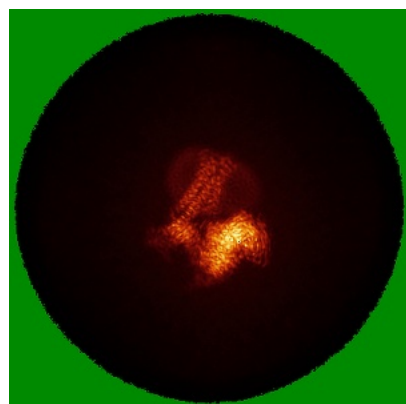


Z Index: 170

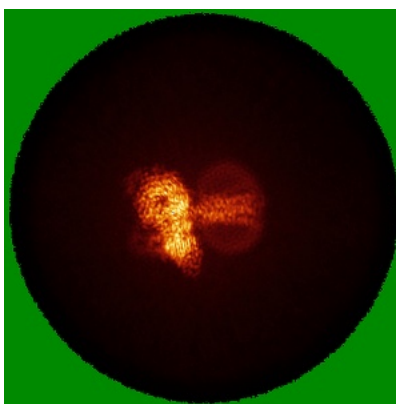
The images above show the largest variance slices of the map in three orthogonal directions.

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

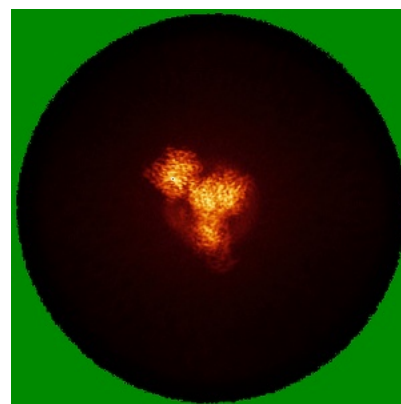
5.4.1 Primary map



X

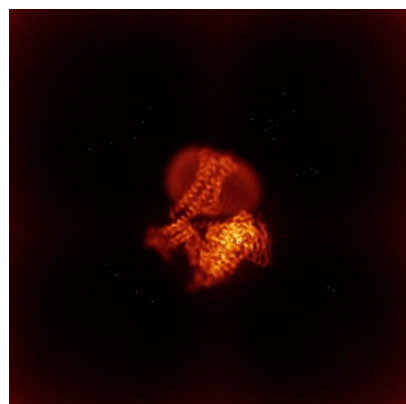


Y

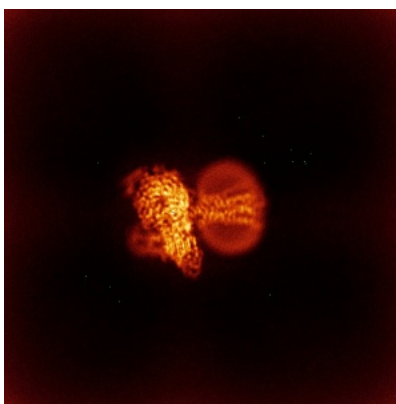


Z

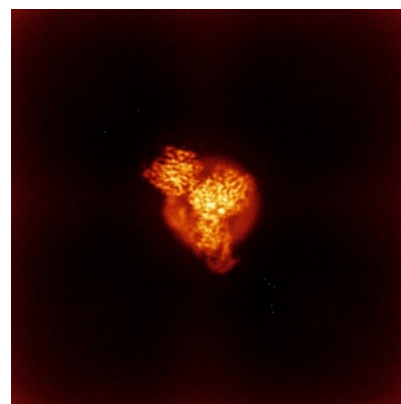
5.4.2 Raw map



X



Y

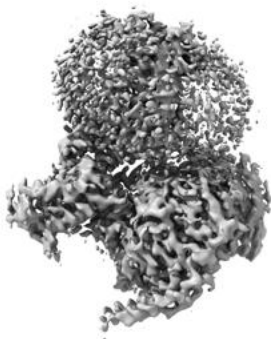


Z

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.

5.5 Orthogonal surface views [i](#)

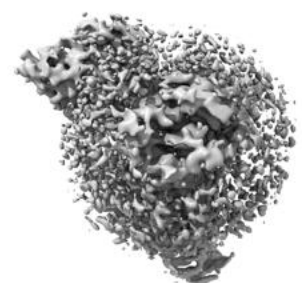
5.5.1 Primary map



X



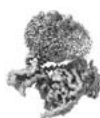
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.12. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

5.5.2 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.

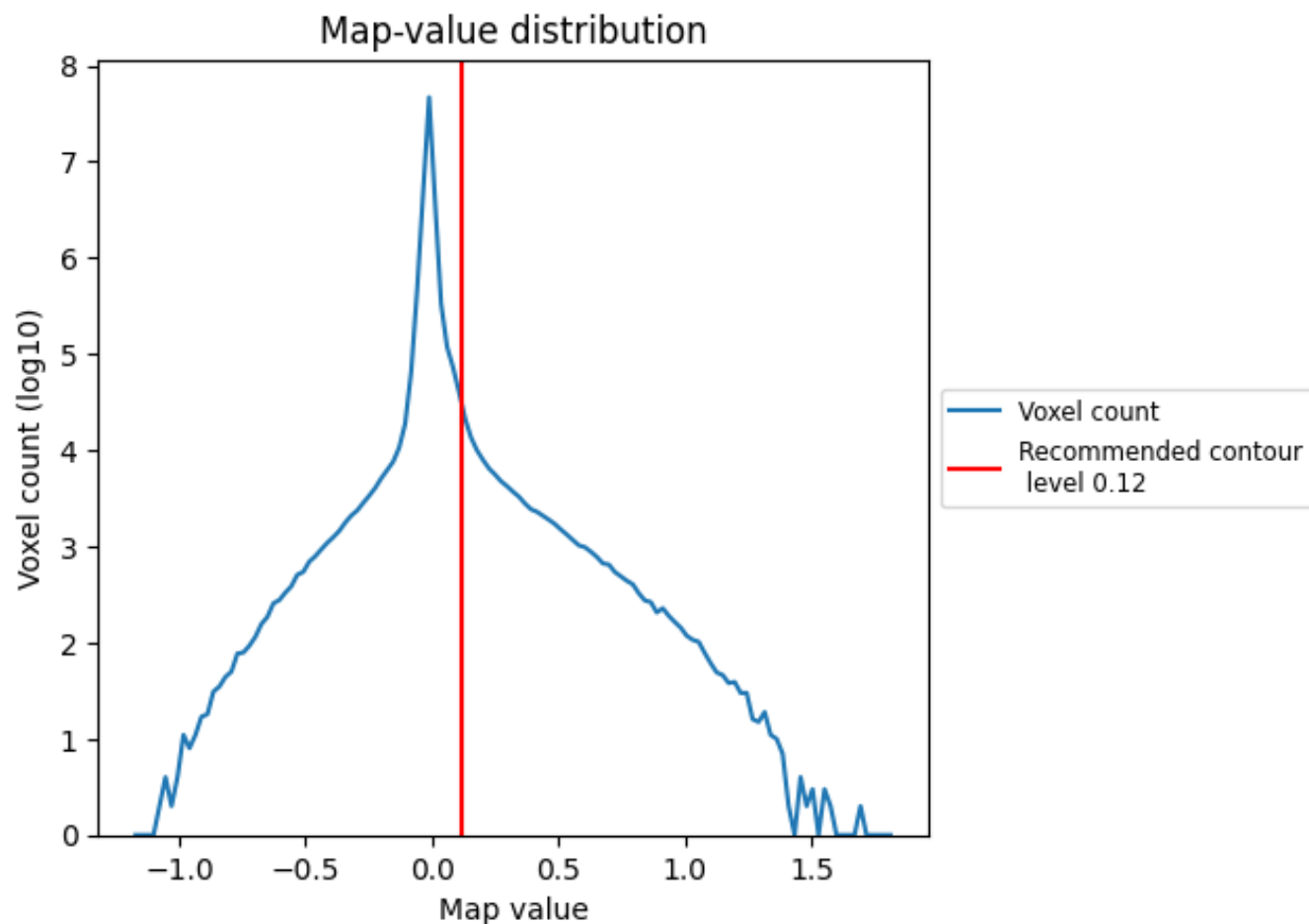
5.6 Mask visualisation [i](#)

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

6 Map analysis [i](#)

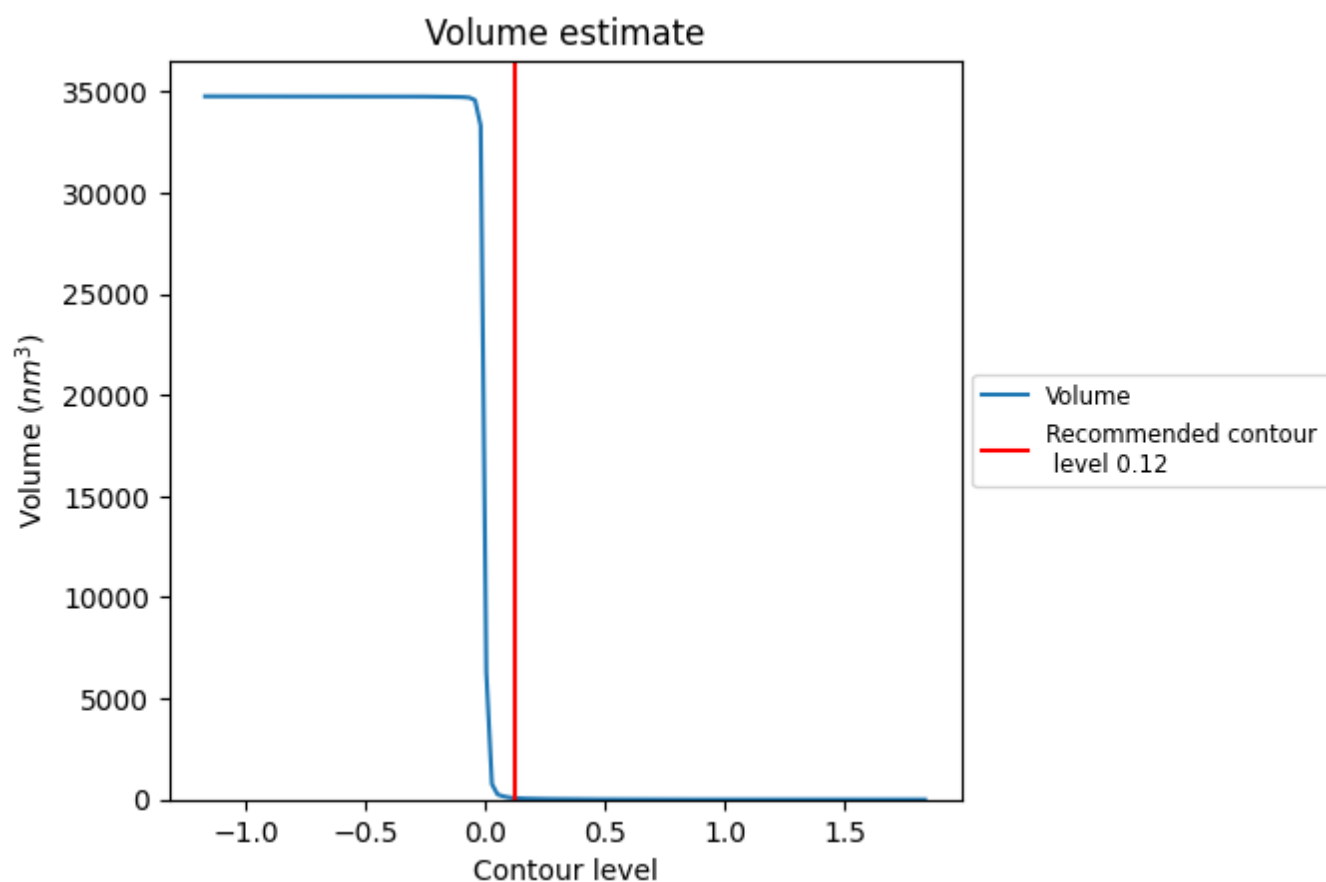
This section contains the results of statistical analysis of the map.

6.1 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.

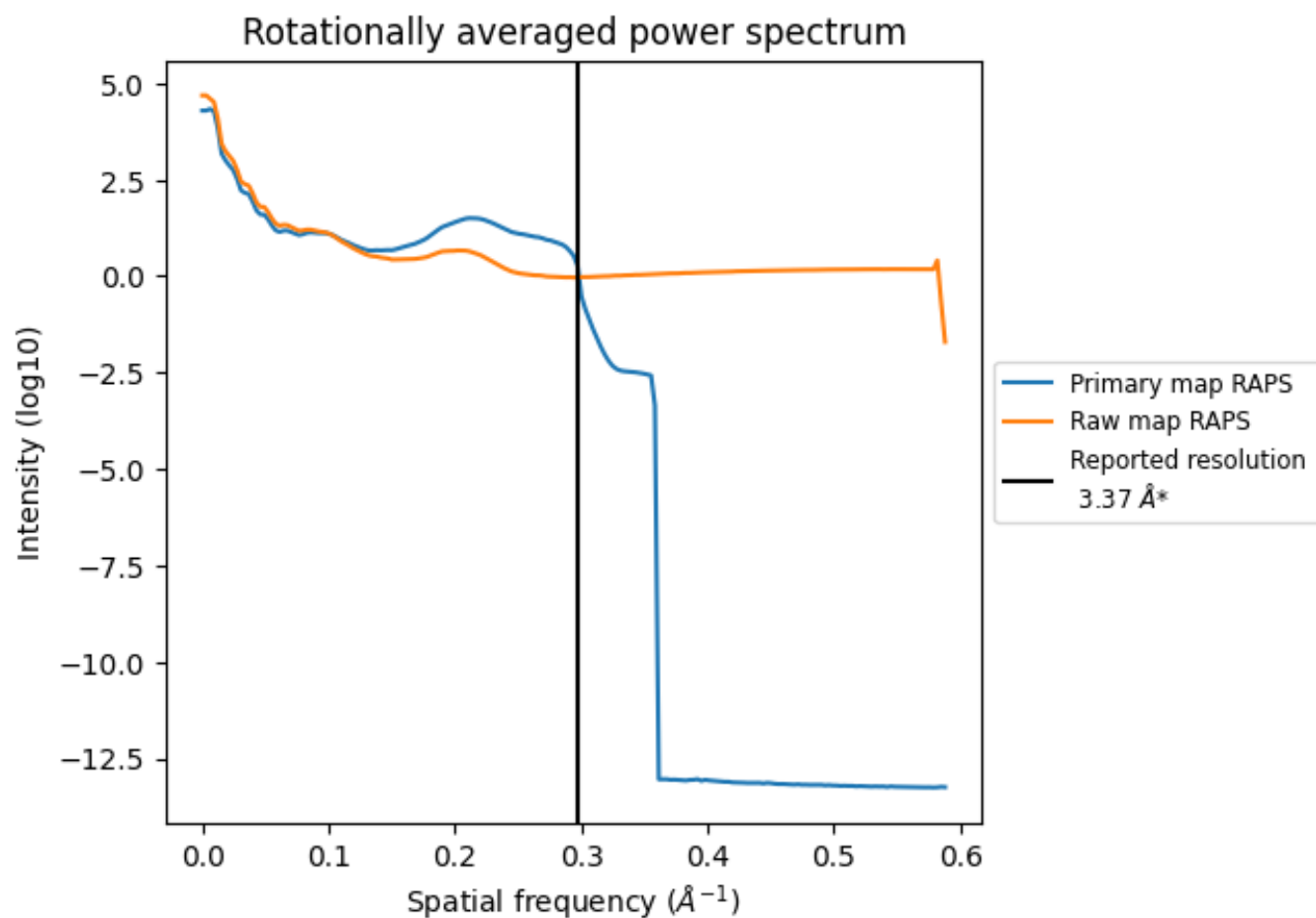
6.2 Volume estimate [i](#)



The volume at the recommended contour level is 75 nm³; this corresponds to an approximate mass of 68 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.

6.3 Rotationally averaged power spectrum ⓘ

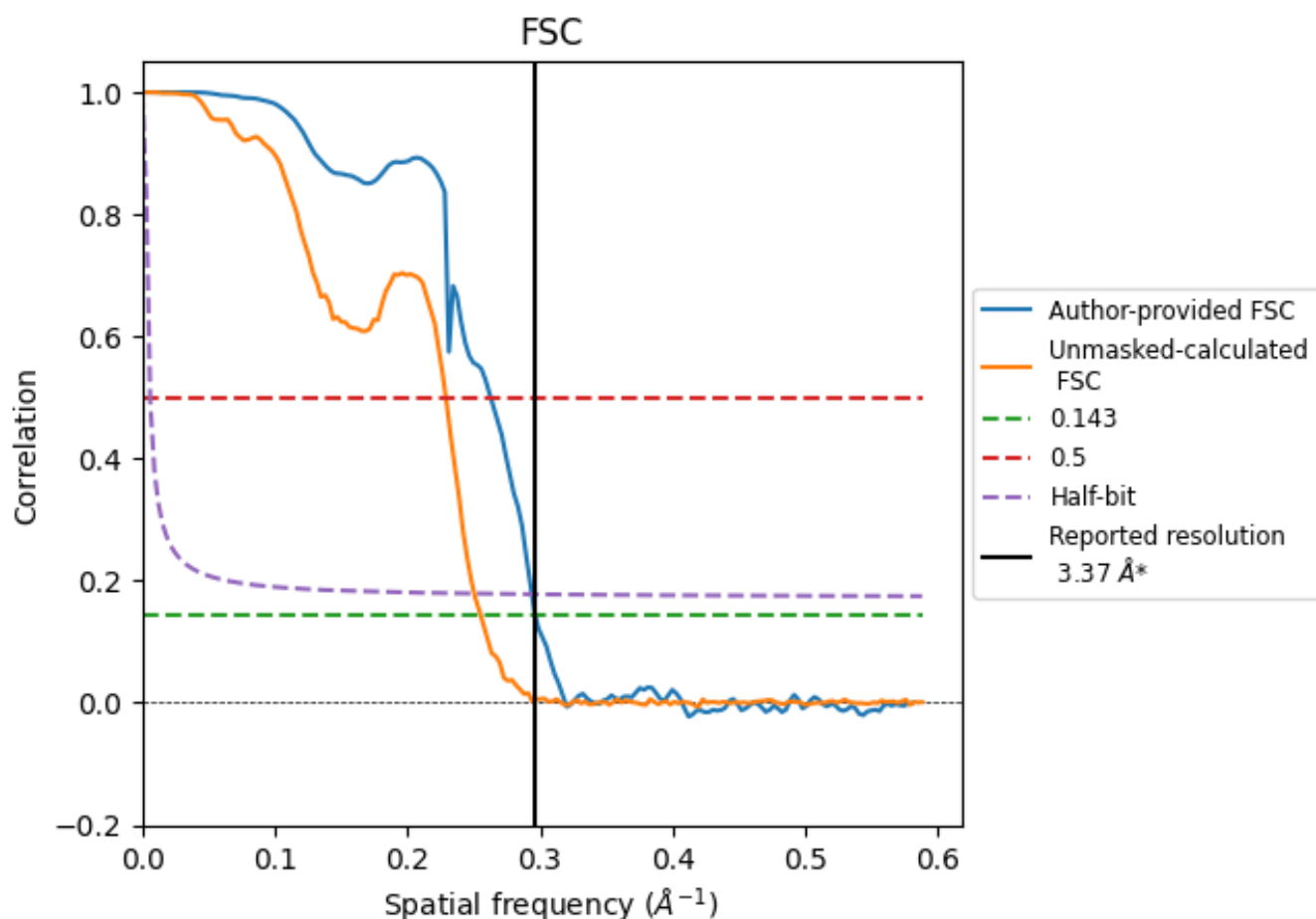


*Reported resolution corresponds to spatial frequency of 0.297 Å⁻¹

7 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.

7.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.297 Å⁻¹

7.2 Resolution estimates [i](#)

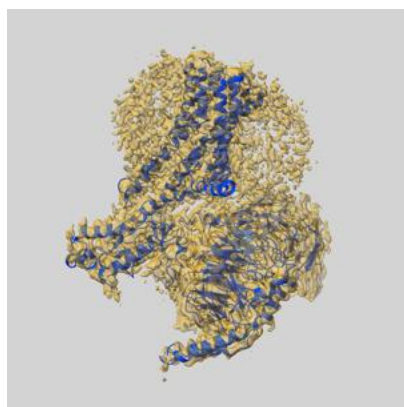
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.37	-	-
Author-provided FSC curve	3.37	3.81	3.40
Unmasked-calculated*	3.92	4.37	3.99

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.92 differs from the reported value 3.37 by more than 10 %

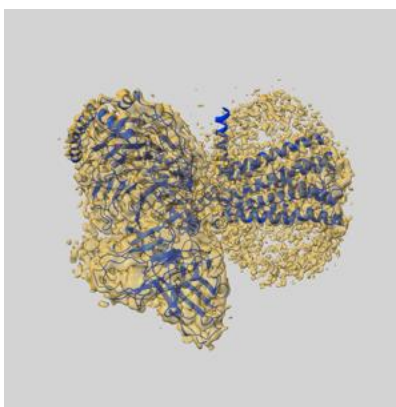
8 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-60651 and PDB model 9IK9. Per-residue inclusion information can be found in section ?? on page ??.

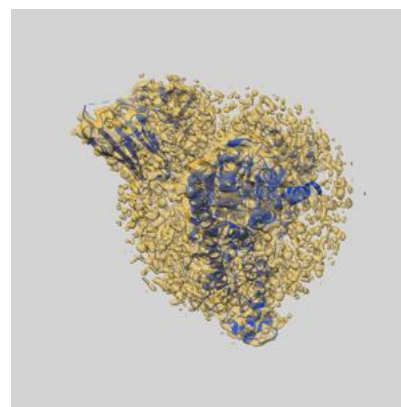
8.1 Map-model overlay [i](#)



X



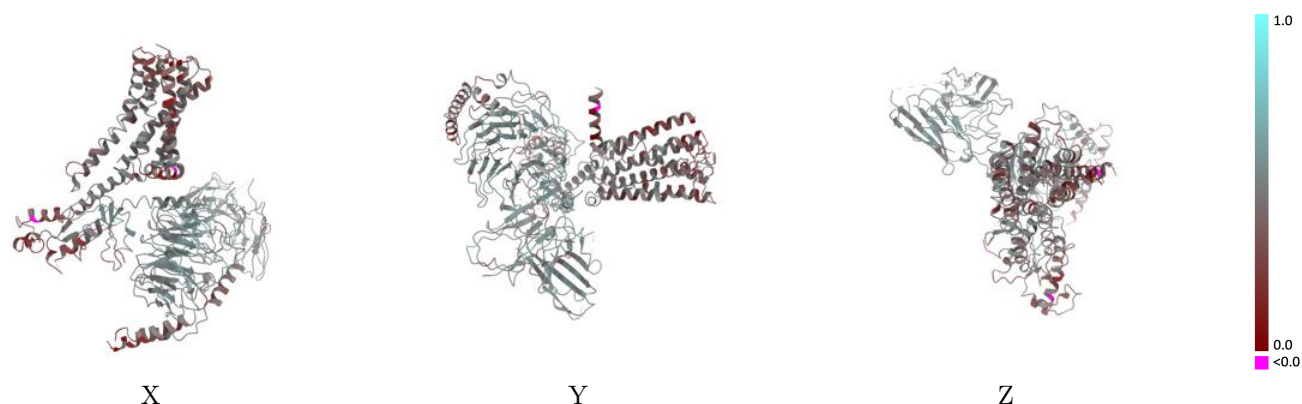
Y



Z

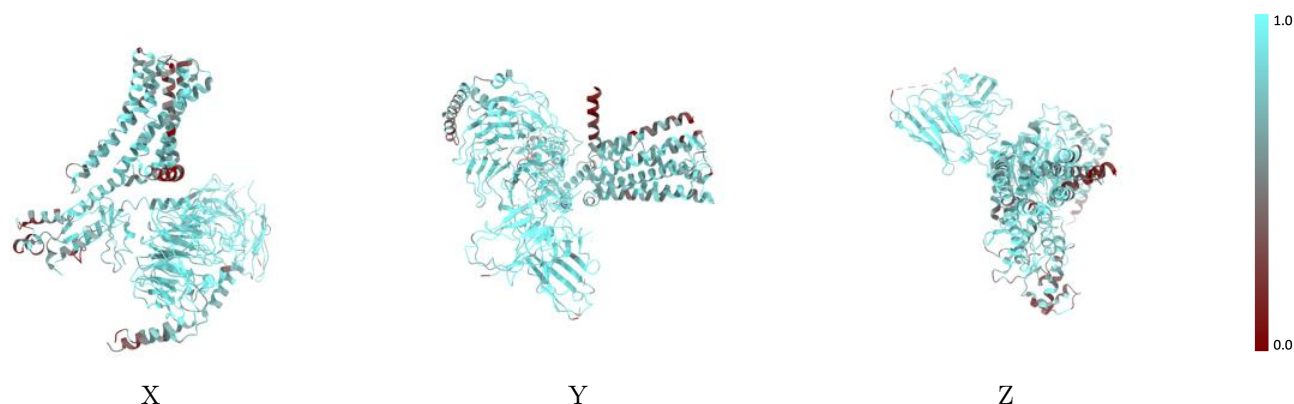
The images above show the 3D surface view of the map at the recommended contour level 0.12 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



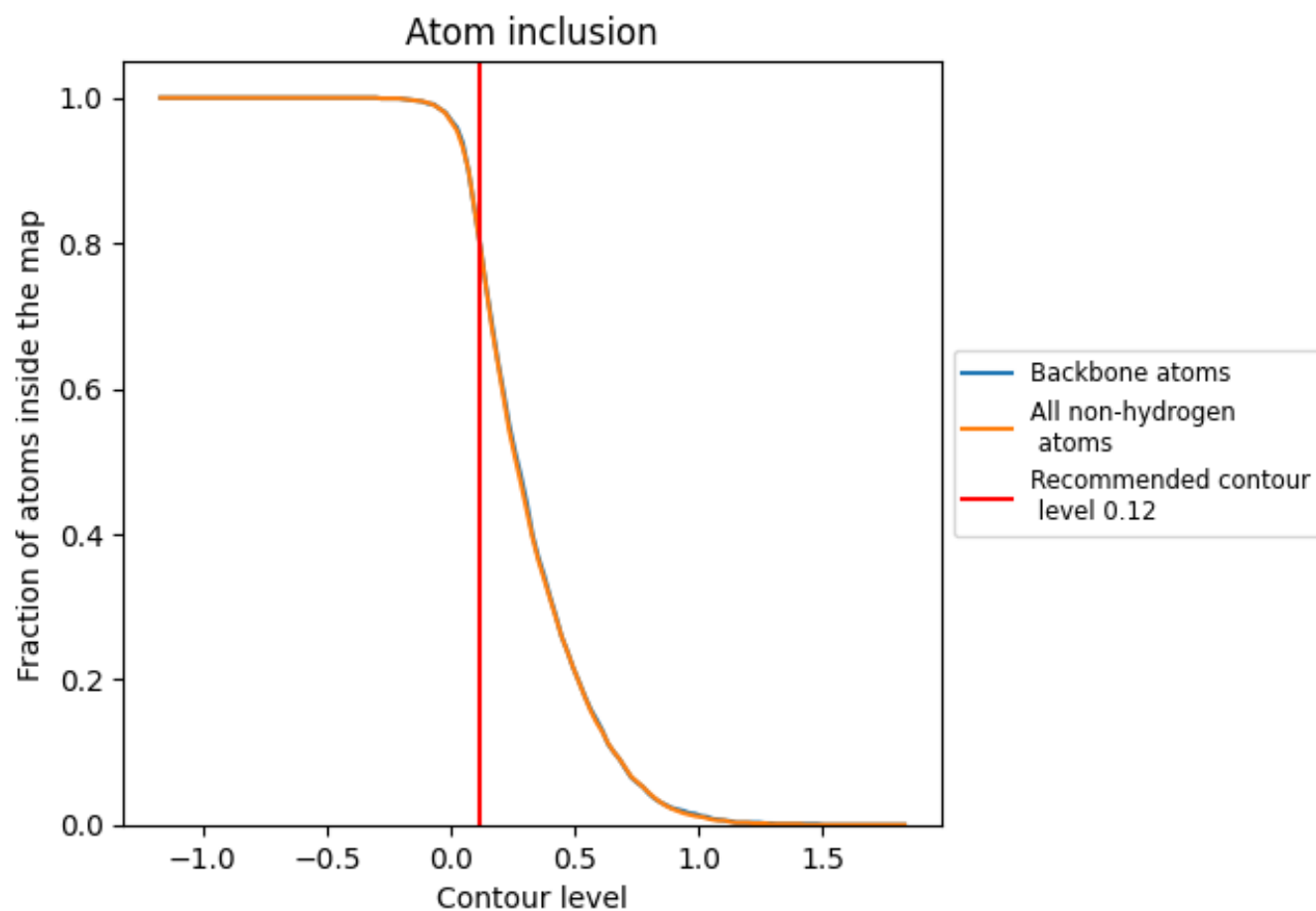
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

8.3 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 (0.12).

8.4 Atom inclusion [i](#)



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

8.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7960	0.4640
A	0.7320	0.4350
B	0.8670	0.5150
C	0.7040	0.4350
D	0.7050	0.3930
E	0.8780	0.5200
F	0.7200	0.3730

1.0

0.0

<0.0