



wwPDB EM Validation Summary Report ⓘ

Jan 16, 2023 – 04:24 pm GMT

PDB ID : 8B0J
EMDB ID : EMD-15785
Title : CryoEM structure of bacterial RNaseE.RapZ.GlmZ complex central to the control of cell envelope biogenesis
Authors : Islam, M.S.; Hardwick, H.W.; Chirgadze, D.Y.; Luisi, B.F.
Deposited on : 2022-09-07
Resolution : 3.99 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:

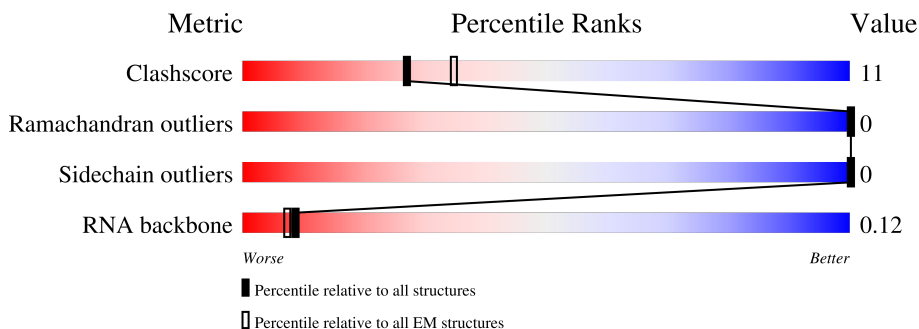
EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.99 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
RNA backbone	4643	859

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	284	
1	B	284	
1	C	284	
1	D	284	
2	L	581	
2	N	581	
3	K	207	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 28636 atoms, of which 12434 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 RNase adapter protein RapZ.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
1	A	268	Total	C	H	N	O	S	0	0
			3695	1219	1780	340	351	5		
1	B	279	Total	C	H	N	O	S	0	0
			3571	1213	1657	340	359	2		
1	D	281	Total	C	H	N	O	S	0	0
			3462	1185	1585	323	366	3		
1	C	266	Total	C	H	N	O	S	0	0
			3239	1115	1484	300	338	2		

- Molecule 2 is a protein called Ribonuclease E.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	L	481	Total	C	H	N	O	S	0	0
			5605	1902	2545	551	606	1		
2	N	356	Total	C	H	N	O	S	0	0
			4220	1428	1915	423	450	4		

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	346	CYS	ASP	conflict	UNP P21513
L	?	-	GLN	deletion	UNP P21513
L	?	-	PRO	deletion	UNP P21513
L	?	-	GLY	deletion	UNP P21513
L	?	-	LEU	deletion	UNP P21513
L	?	-	LEU	deletion	UNP P21513
L	?	-	SER	deletion	UNP P21513
L	?	-	ARG	deletion	UNP P21513
L	?	-	PHE	deletion	UNP P21513
L	?	-	PHE	deletion	UNP P21513
L	?	-	GLY	deletion	UNP P21513
L	?	-	ALA	deletion	UNP P21513
L	?	-	LEU	deletion	UNP P21513

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Chain	Residue	Modelled	Actual	Comment	Reference
L	?	-	LYS	deletion	UNP P21513
L	?	-	ALA	deletion	UNP P21513
L	?	-	LEU	deletion	UNP P21513
L	?	-	PHE	deletion	UNP P21513
L	?	-	SER	deletion	UNP P21513
N	346	CYS	ASP	conflict	UNP P21513
N	?	-	GLN	deletion	UNP P21513
N	?	-	PRO	deletion	UNP P21513
N	?	-	GLY	deletion	UNP P21513
N	?	-	LEU	deletion	UNP P21513
N	?	-	LEU	deletion	UNP P21513
N	?	-	SER	deletion	UNP P21513
N	?	-	ARG	deletion	UNP P21513
N	?	-	PHE	deletion	UNP P21513
N	?	-	PHE	deletion	UNP P21513
N	?	-	GLY	deletion	UNP P21513
N	?	-	ALA	deletion	UNP P21513
N	?	-	LEU	deletion	UNP P21513
N	?	-	LYS	deletion	UNP P21513
N	?	-	ALA	deletion	UNP P21513
N	?	-	LEU	deletion	UNP P21513
N	?	-	PHE	deletion	UNP P21513
N	?	-	SER	deletion	UNP P21513

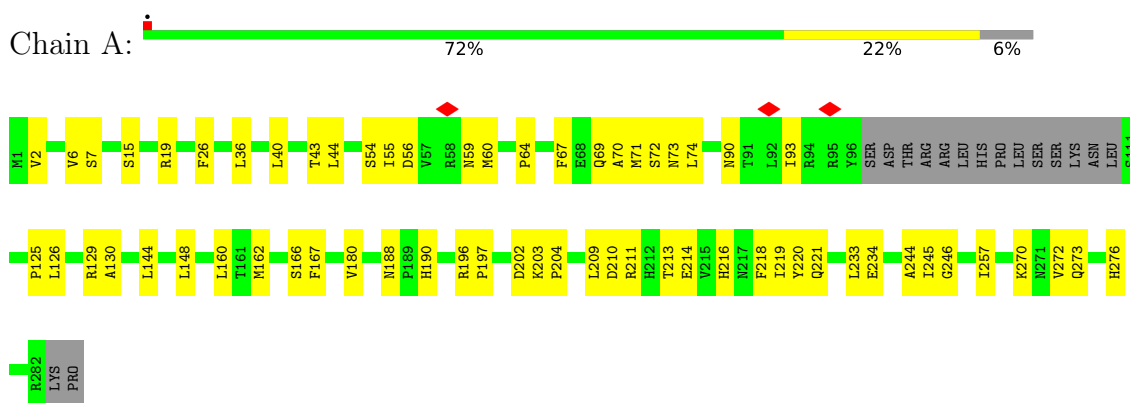
- Molecule 3 is a RNA chain called GlmZ small RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
3	K	160	4844	1511	1468	576	1129	160	0	0

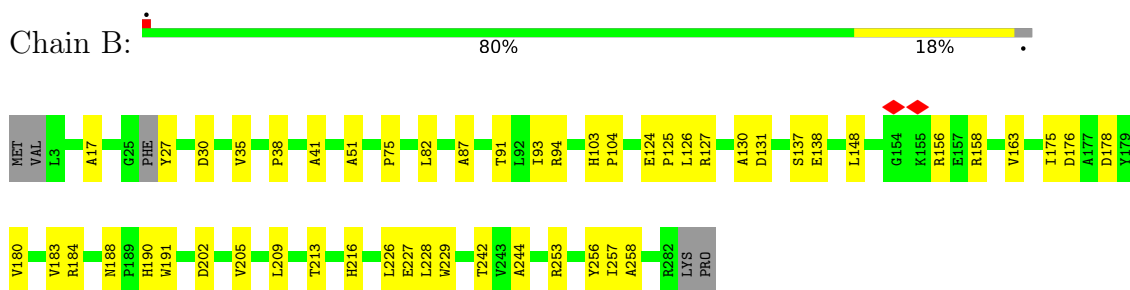
3 Residue-property plots [i](#)

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.

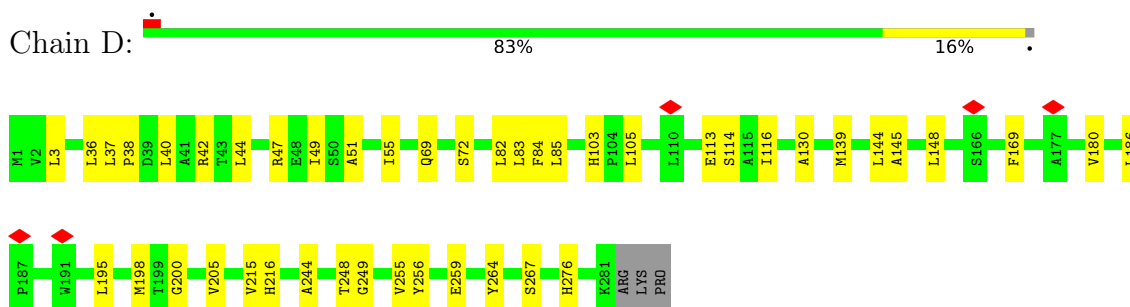
- Molecule 1: RNase adapter protein RapZ



- Molecule 1: RNase adapter protein RapZ



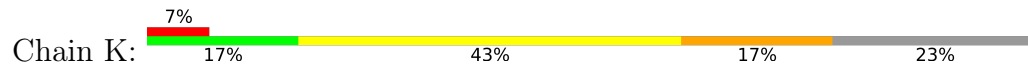
- Molecule 1: RNase adapter protein RapZ



- Molecule 1: RNase adapter protein RapZ

PRO GLU GLN PRO PRO ALA LEU ALA THR PHE MET PRO ASP VAL PRO PRO ALA THR PRO ALA VAL VAL PRO ALA ALA PRO PRO LYS ALA ALA PRO PRO ALA GLY GLU THR ALA ALA PRO PRO ALA GLY GLU THR LYS PRO THR GLN PRO ALA LYS ALA

● Molecule 3: GlmZ small RNA



G U A G A U G7 C8 U9 C10 A11 U12 U13 C14 C15 A16 U17 C18 U19 C20 U21 U22 A23 U24 U27 C28 G29 C30 C31 U32 U33 A34 G35 U36 G37 U40 C41 A42 U43 A44 A45 A46 C47 U48 C49 C50 G51 G52 A53 A54 U55 U56 A57 C58 G59 C60 A61

G64 C C G U U A C G73 U74 U75 G76 C77 U78 U79 A80 U81 C82 C83 U84 C85 A87 C88 U89 G90 A91 C92 A93 G A U G97 U98 C99 G100 C101 U102 U103 A U G U C U109 C110 A111 G115 A C A C A121 U122 G G124 A125

C126 A127 C128 A129 C130 G131 G132 U133 U134 G135 A136 G137 U138 G139 A144 C145 C146 C147 A148 C149 U150 U151 G152 U U G U C A U A C A G A C U U G U U U U172 G176 C177 C178 U179 G180 C181 U182 C183 C184 G185 U186 U187 A188 A189 U190

A191 A192 G193 C196 A197 G198 G199 C200 G201 U202 U203 U204 U205 U206 U207

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	33595	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	47.30	Depositor
Minimum defocus (nm)	100	Depositor
Maximum defocus (nm)	300	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.309	Depositor
Minimum map value	-0.193	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.04	Depositor
Map size (\AA)	299.92, 299.92, 299.92	wwPDB
Map dimensions	460, 460, 460	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.652, 0.652, 0.652	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	0.59	0/1951	0.61	0/2666
1	B	0.59	0/1952	0.61	0/2679
1	C	0.49	0/1785	0.58	0/2460
1	D	0.48	0/1909	0.61	0/2622
2	L	0.50	0/3094	0.63	0/4250
2	N	0.47	0/2325	0.61	0/3179
3	K	0.38	1/3760 (0.0%)	0.89	0/5835
All	All	0.49	1/16776 (0.0%)	0.69	0/23691

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	85	C	O3'-P	5.94	1.68	1.61

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1915	1780	1781	40	0
1	B	1914	1657	1655	35	0
1	C	1755	1484	1483	30	0
1	D	1877	1585	1585	32	0
2	L	3060	2545	2542	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	N	2305	1915	1919	32	0
3	K	3376	1468	1721	107	0
All	All	16202	12434	12686	314	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.

The worst 5 of 314 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)
3:K:58:C:N4	3:K:59:G:C6	2.04	1.25
3:K:56:G:O2'	3:K:57:A:H5'	1.39	1.20
3:K:179:U:OP2	3:K:180:G:C6	1.95	1.19
3:K:58:C:C4	3:K:59:G:C5	2.37	1.11
3:K:131:C:H1'	3:K:132:G:H5'	1.28	1.10

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	264/284 (93%)	244 (92%)	20 (8%)	0	100	100
1	B	275/284 (97%)	242 (88%)	33 (12%)	0	100	100
1	C	262/284 (92%)	242 (92%)	20 (8%)	0	100	100
1	D	279/284 (98%)	250 (90%)	29 (10%)	0	100	100
2	L	473/581 (81%)	428 (90%)	45 (10%)	0	100	100
2	N	340/581 (58%)	303 (89%)	37 (11%)	0	100	100
All	All	1893/2298 (82%)	1709 (90%)	184 (10%)	0	100	100

There are no Ramachandran outliers to report.

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	179/254 (70%)	179 (100%)	0	100	100
1	B	161/254 (63%)	161 (100%)	0	100	100
1	C	142/254 (56%)	142 (100%)	0	100	100
1	D	152/254 (60%)	152 (100%)	0	100	100
2	L	227/484 (47%)	227 (100%)	0	100	100
2	N	174/484 (36%)	174 (100%)	0	100	100
All	All	1035/1984 (52%)	1035 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	268	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	K	156/207 (75%)	97 (62%)	24 (15%)

5 of 97 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	K	9	U
3	K	10	C
3	K	14	C
3	K	16	A
3	K	17	U

5 of 24 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	K	138	U
3	K	150	U
3	K	149	C
3	K	176	G
3	K	97	G

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](#)

There are no chain breaks in this entry.

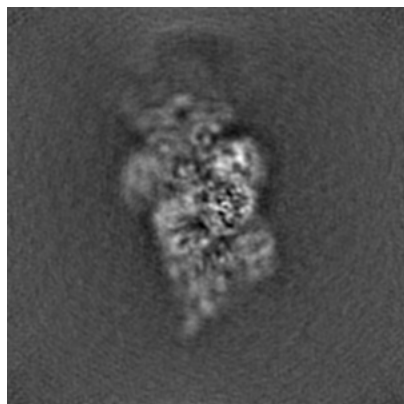
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15785. These allow visual inspection of the internal detail of the map and identification of artifacts.

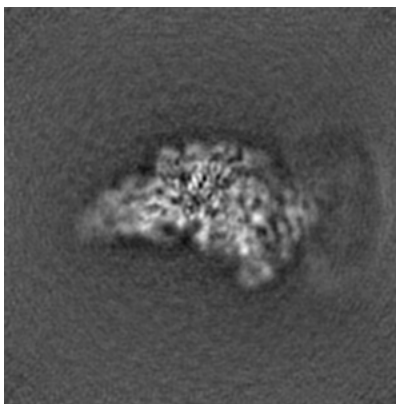
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

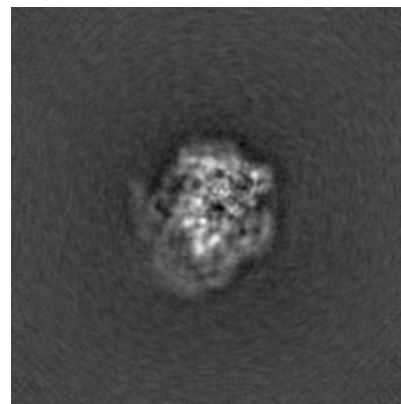
6.1.1 Primary map



X

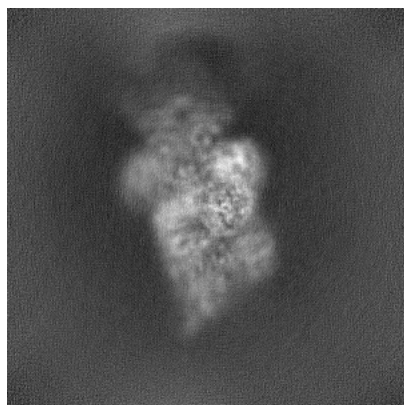


Y

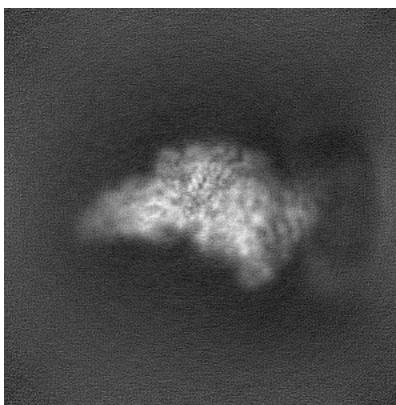


Z

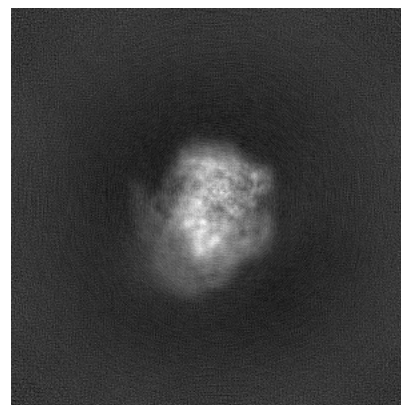
6.1.2 Raw map



X



Y

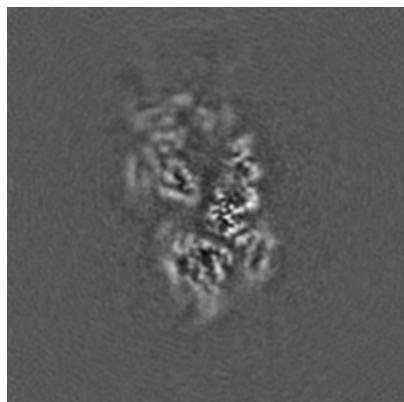


Z

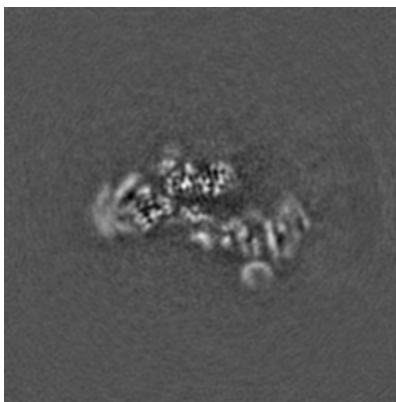
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

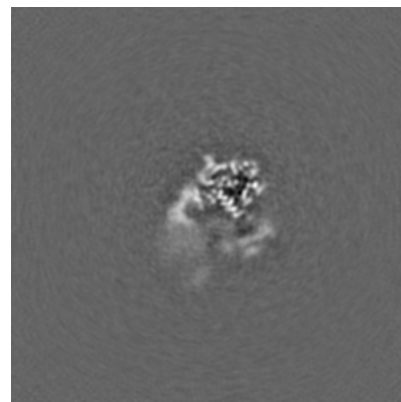
6.2.1 Primary map



X Index: 230

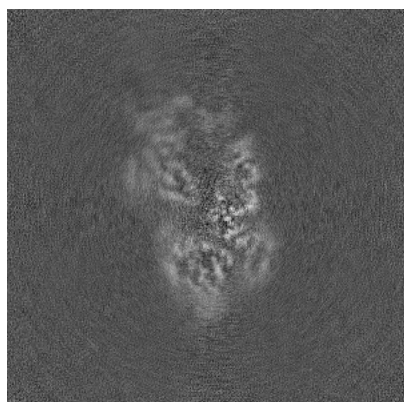


Y Index: 230

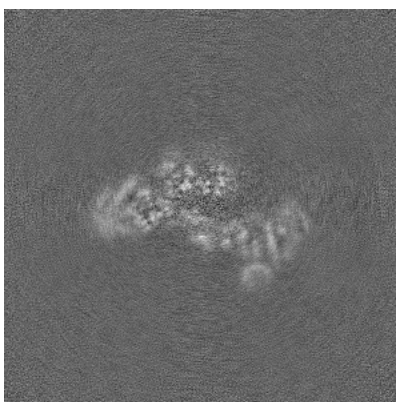


Z Index: 230

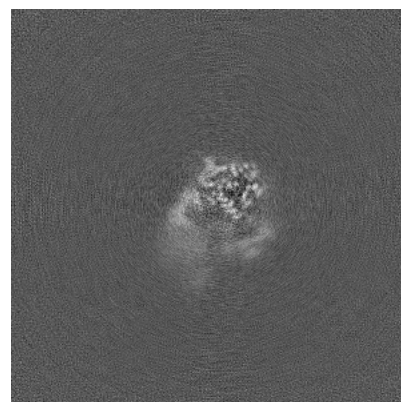
6.2.2 Raw map



X Index: 230



Y Index: 230

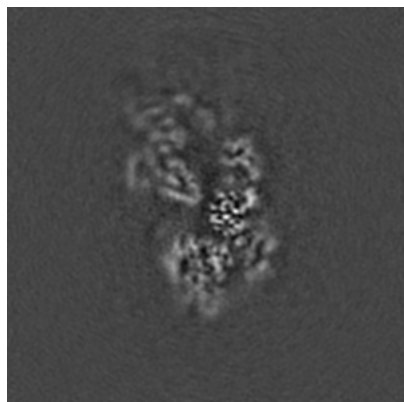


Z Index: 230

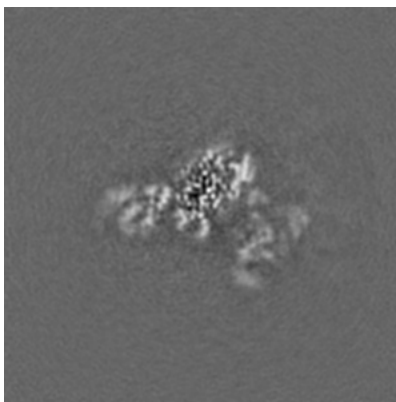
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

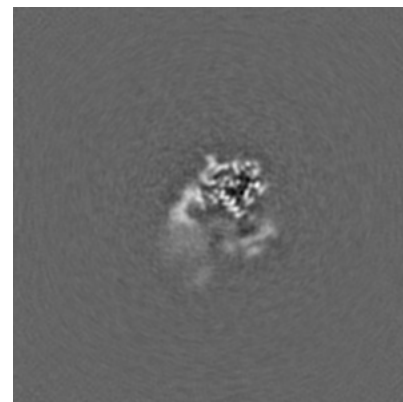
6.3.1 Primary map



X Index: 233

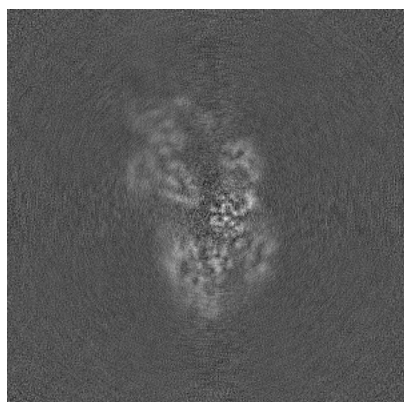


Y Index: 247

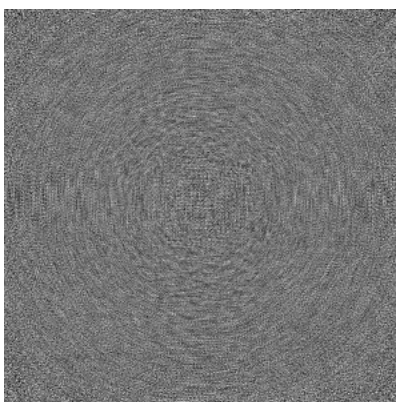


Z Index: 230

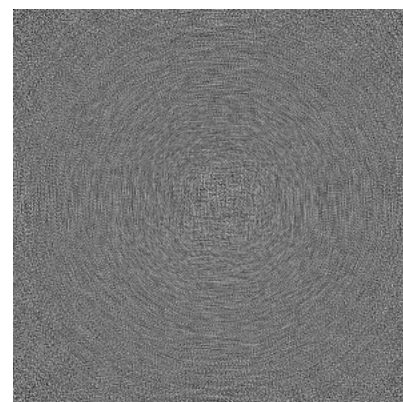
6.3.2 Raw map



X Index: 233



Y Index: 0

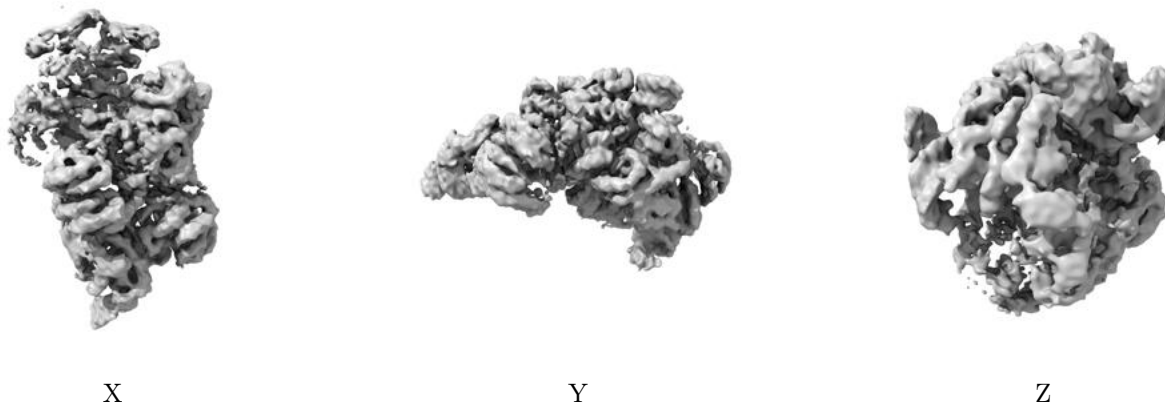


Z Index: 0

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

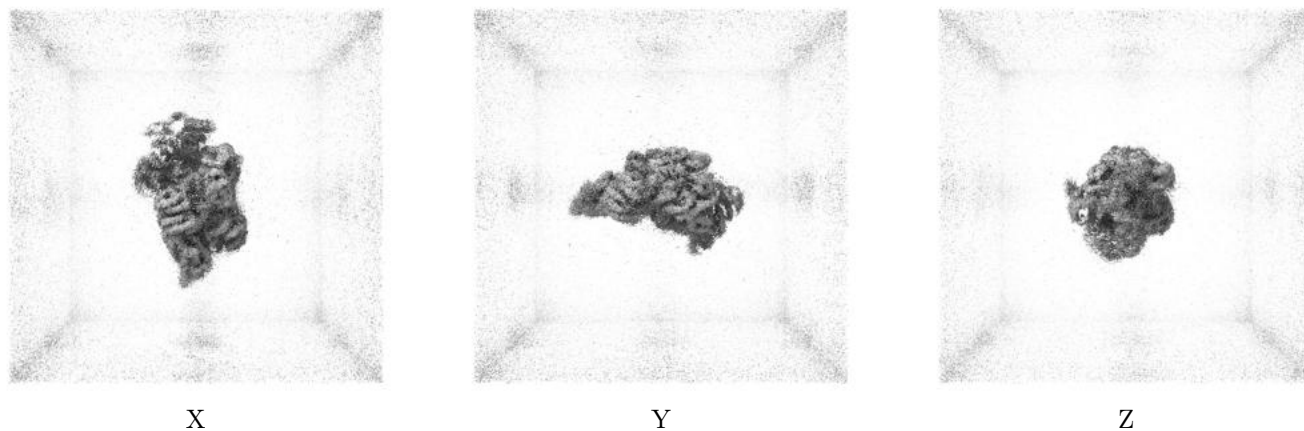
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.04. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

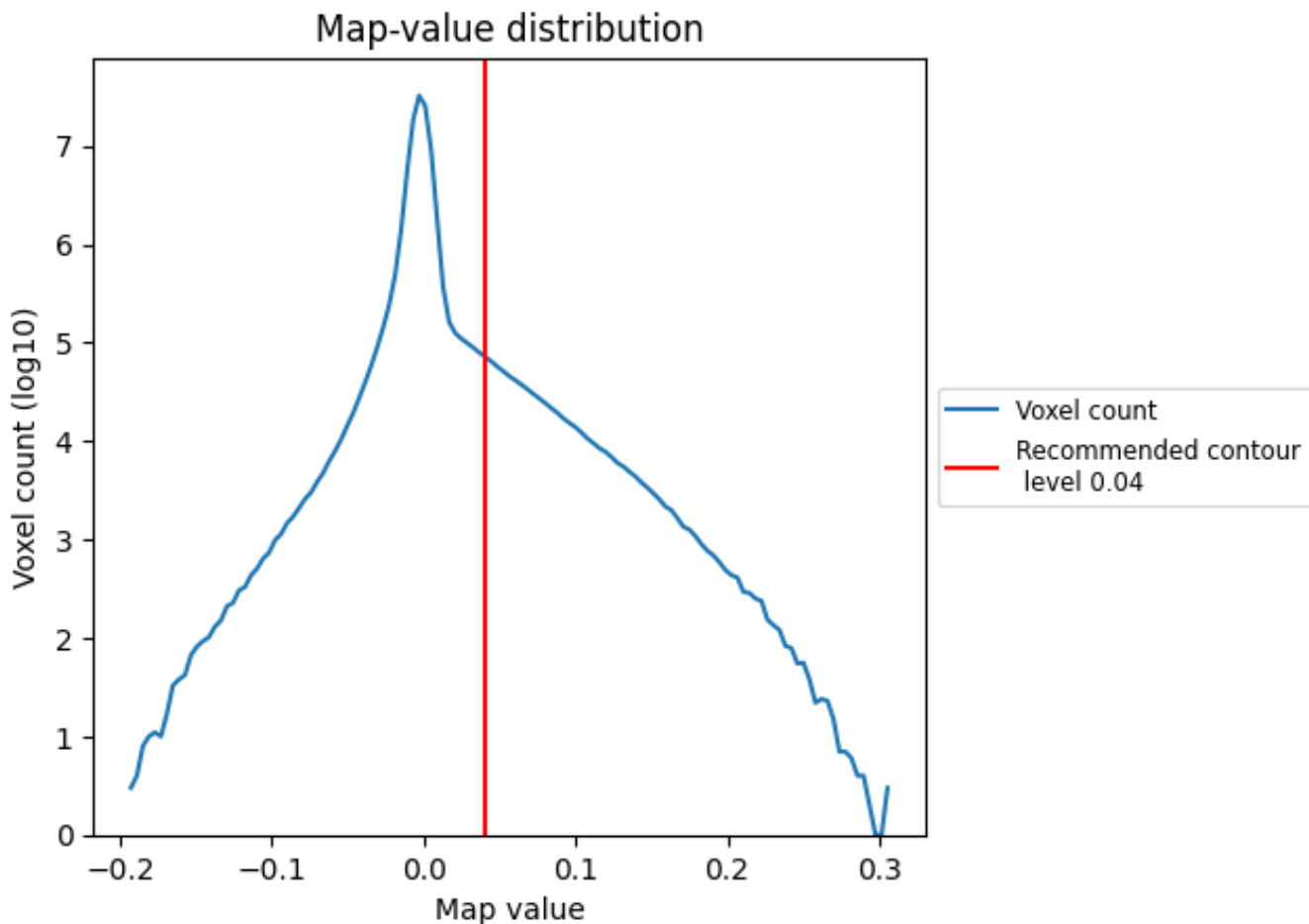
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

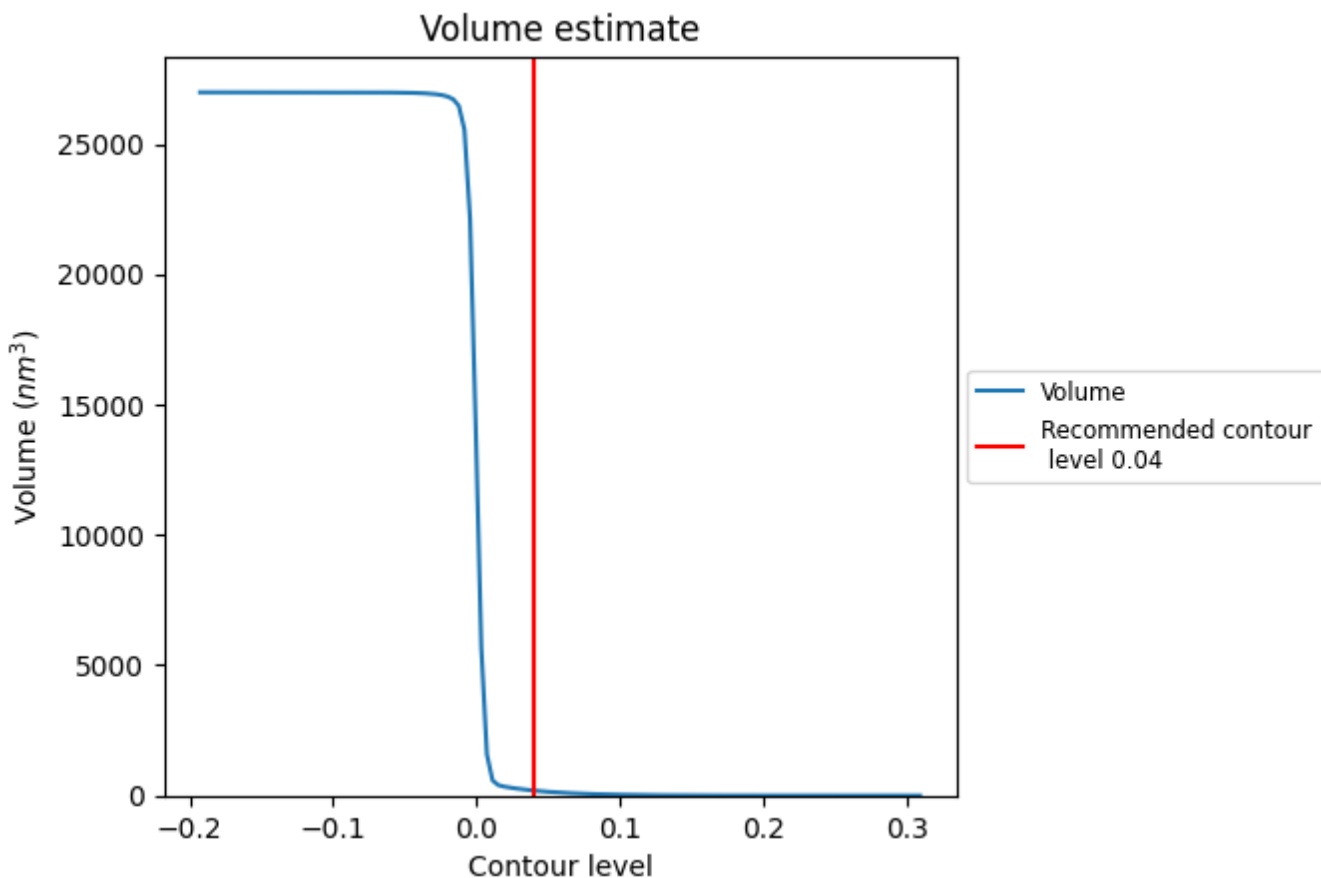
This section contains the results of statistical analysis of the map.

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.

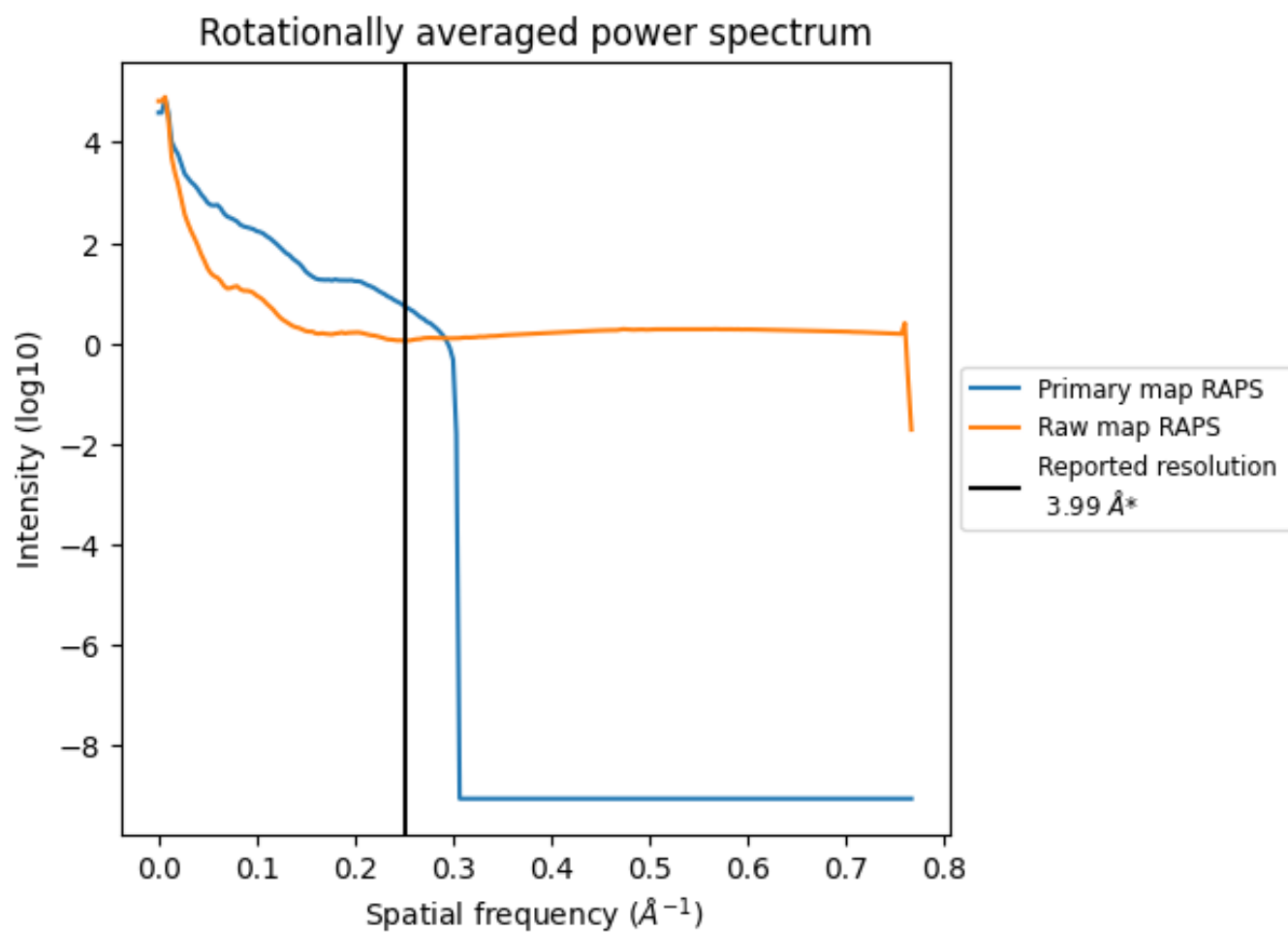
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 191 nm^3 ; this corresponds to an approximate mass of 173 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](#)

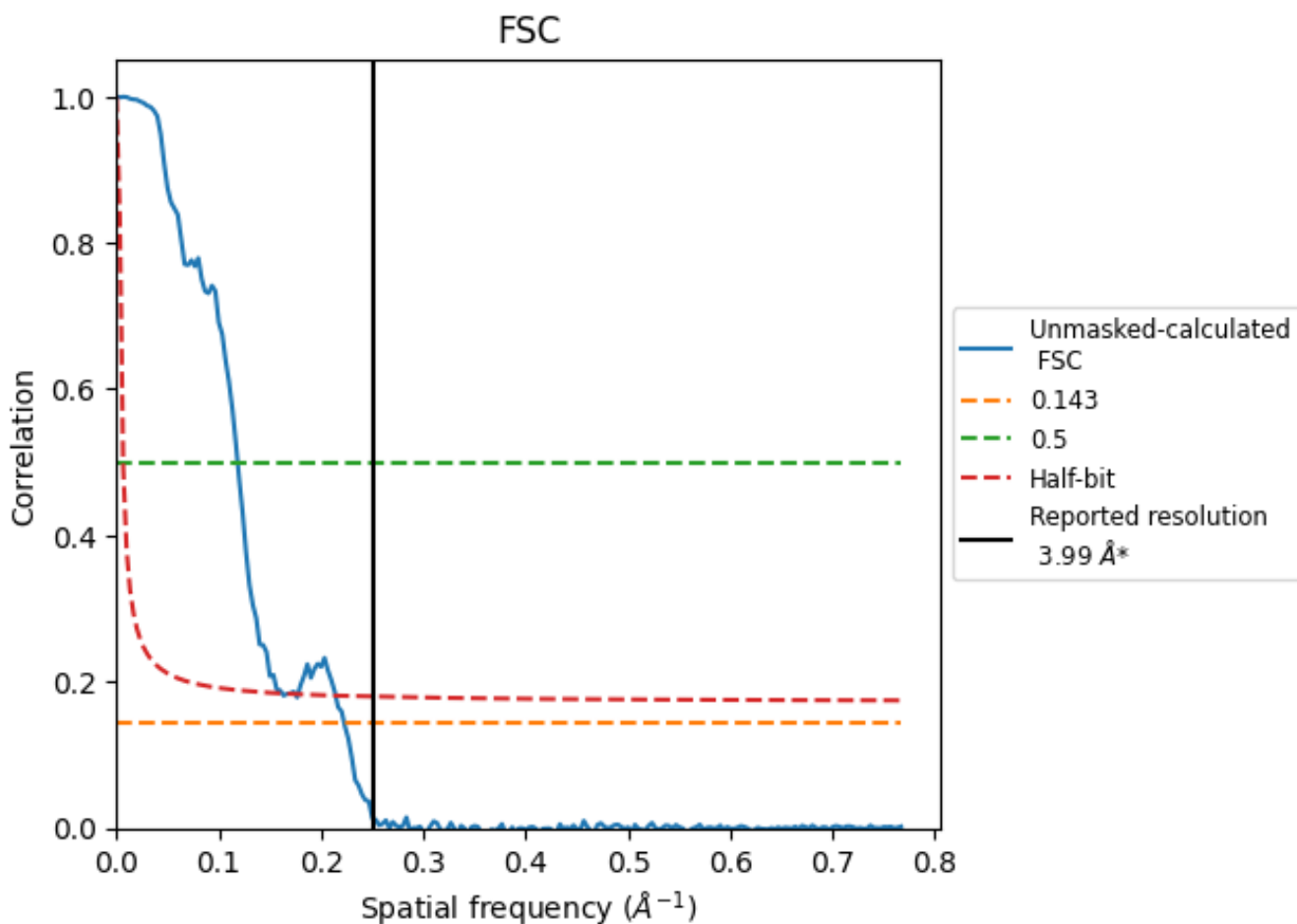


*Reported resolution corresponds to spatial frequency of 0.251 Å⁻¹

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 0.251 Å⁻¹

8.2 Resolution estimates [i](#)

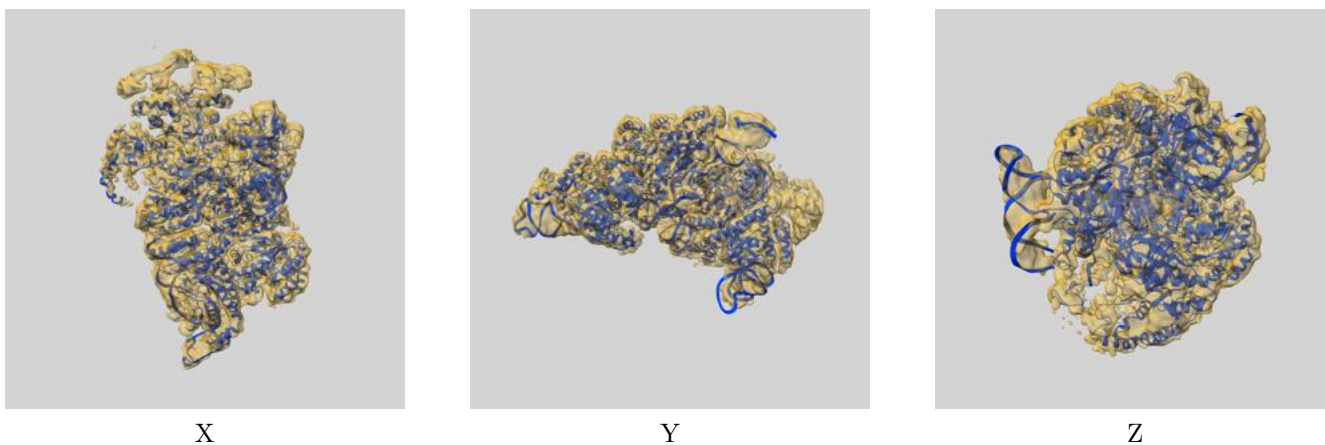
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.99	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.49	8.44	6.17

*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 4.49 differs from the reported value 3.99 by more than 10 %

9 Map-model fit [i](#)

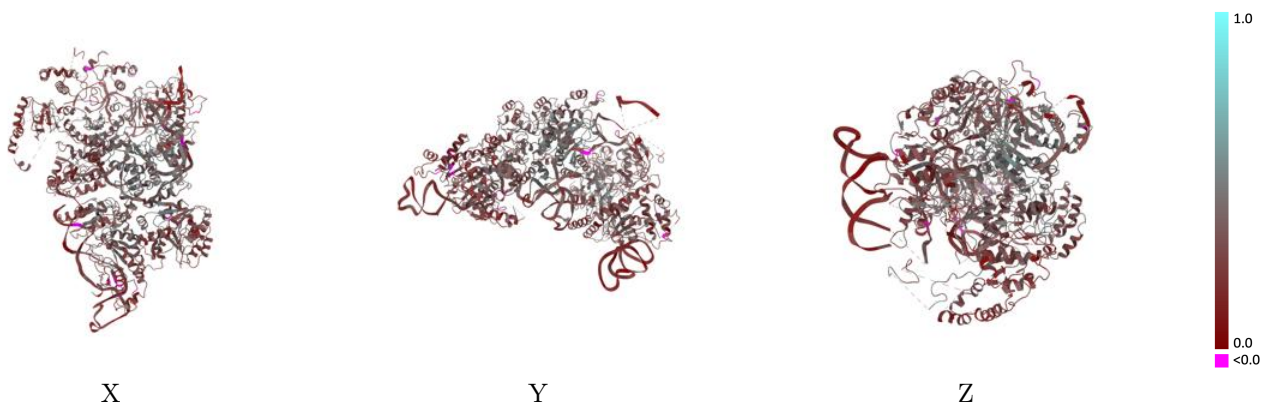
This section contains information regarding the fit between EMDB map EMD-15785 and PDB model 8B0J. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)



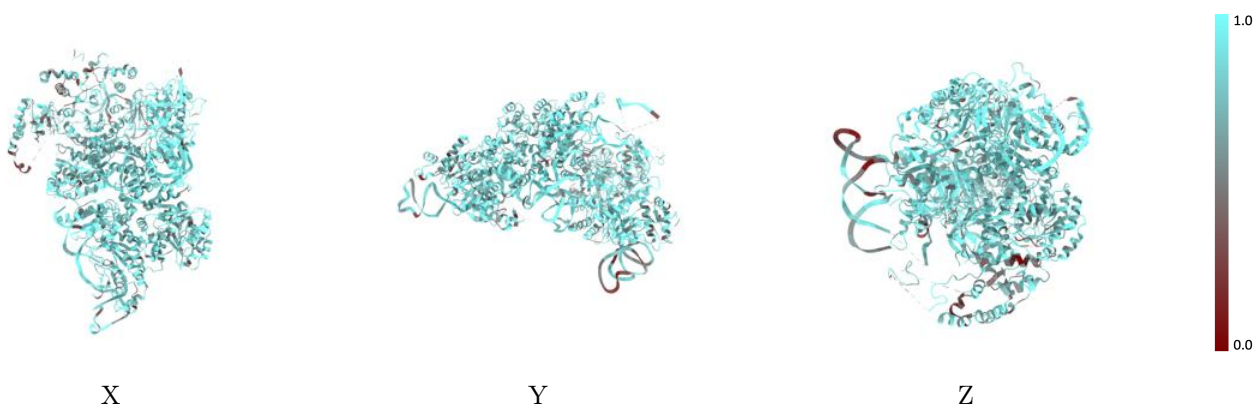
The images above show the 3D surface view of the map at the recommended contour level 0.04 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



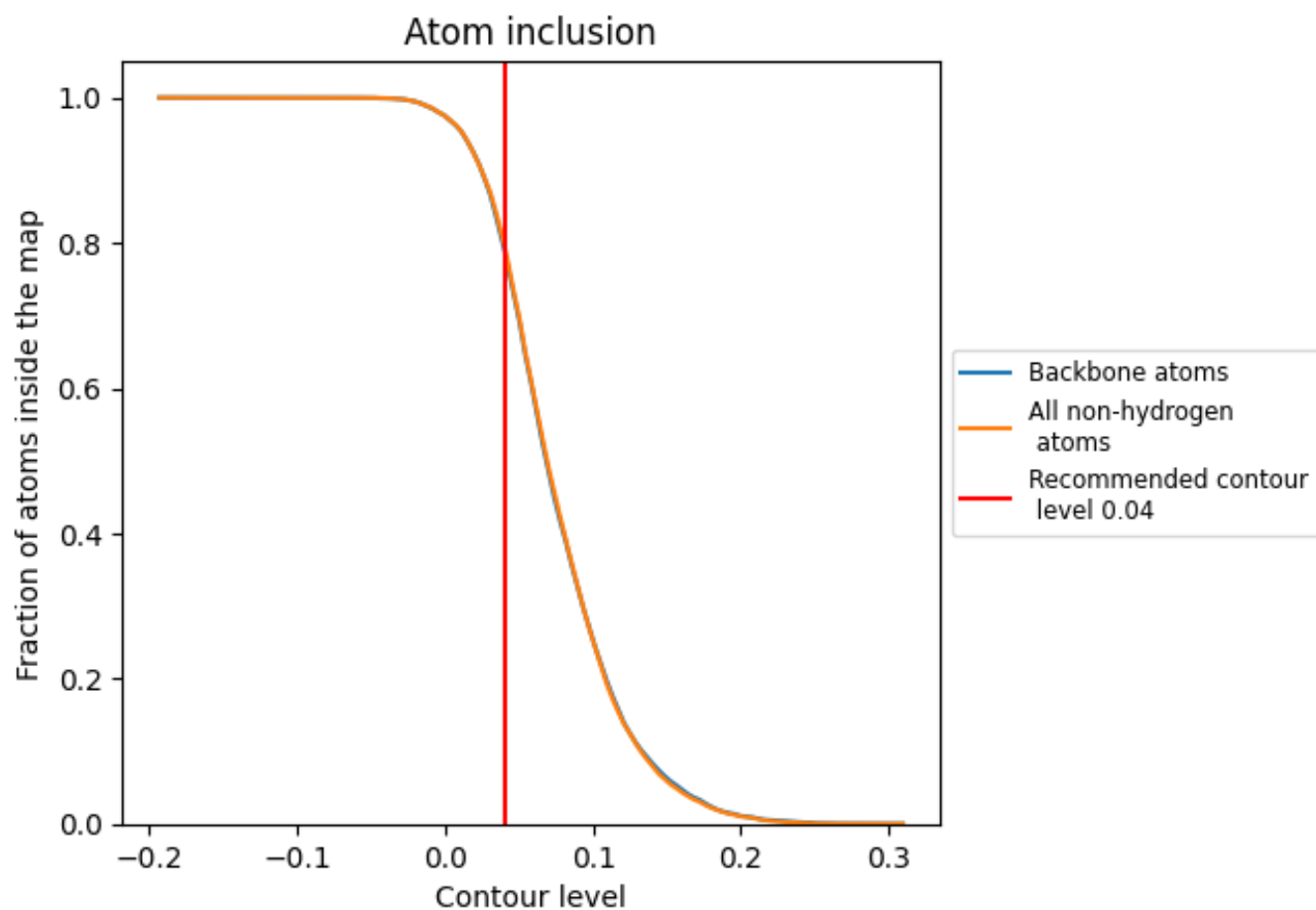
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.04).

















9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 80% 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 (0.04) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7979	 0.3190
A	 0.8307	 0.4020
B	 0.8613	 0.3900
C	 0.8056	 0.3180
D	 0.8146	 0.3160
K	 0.7974	 0.2350
L	 0.8092	 0.3310
N	 0.7214	 0.3030

