



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 11:23 PM UTC

PDB ID : 5TBY / pdb_00005tby
EMDB ID : EMD-2240
Title : HUMAN BETA CARDIAC HEAVY MEROMYOSIN INTERACTING-HEADS MOTIF OBTAINED BY HOMOLOGY MODELING (USING SWISS-MODEL) OF HUMAN SEQUENCE FROM APHONOPELMA HOMOLOGY MODEL (PDB-3JBH), RIGIDLY FITTED TO HUMAN BETA-CARDIAC NEGATIVELY STAINED THICK FILAMENT 3D-RECONSTRUCTION (EMD-2240)
Authors : ALAMO, L.; WARE, J.S.; PINTO, A.; GILLILAN, R.E.; SEIDMAN, J.G.; SEIDMAN, C.E.; PADRON, R.
Deposited on : 2016-09-13
Resolution : 20.00 Å(reported)
Based on initial model : 3JBH

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:

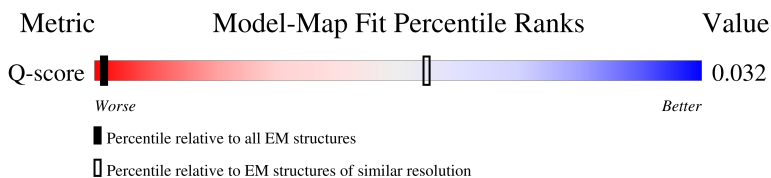
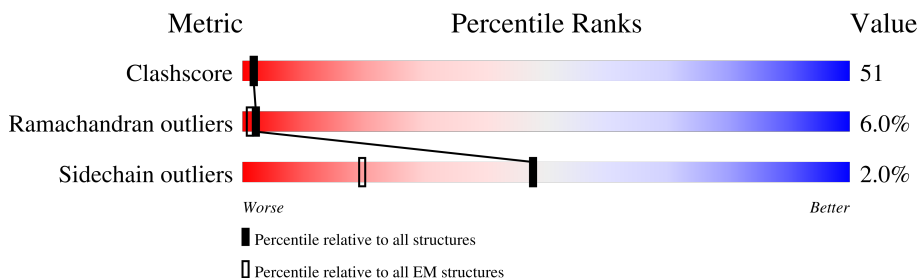
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13

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 20.00 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	11 (27.70 - 28.40)

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	1935	
1	B	1935	

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Ideal geometry (proteins) : Engh & Huber (2001)
 Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
2	C	195	 31% 37% 37% 2% 2% 2%
2	D	195	 41% 38% 35% 2% 2% 2%
3	E	166	 7% 40% 45% 8% 2% 2%
3	F	166	 31% 42% 46% 7% 2% 2%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20357 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 Myosin-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	954	Total	C	N	O	S	0	0
			7704	4899	1324	1439	42		
1	B	950	Total	C	N	O	S	0	0
			7673	4877	1320	1435	41		

- Molecule 2 is a protein called Myosin light chain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	152	Total	C	N	O	S	0	0
			1212	759	202	240	11		
2	D	152	Total	C	N	O	S	0	0
			1212	759	202	240	11		

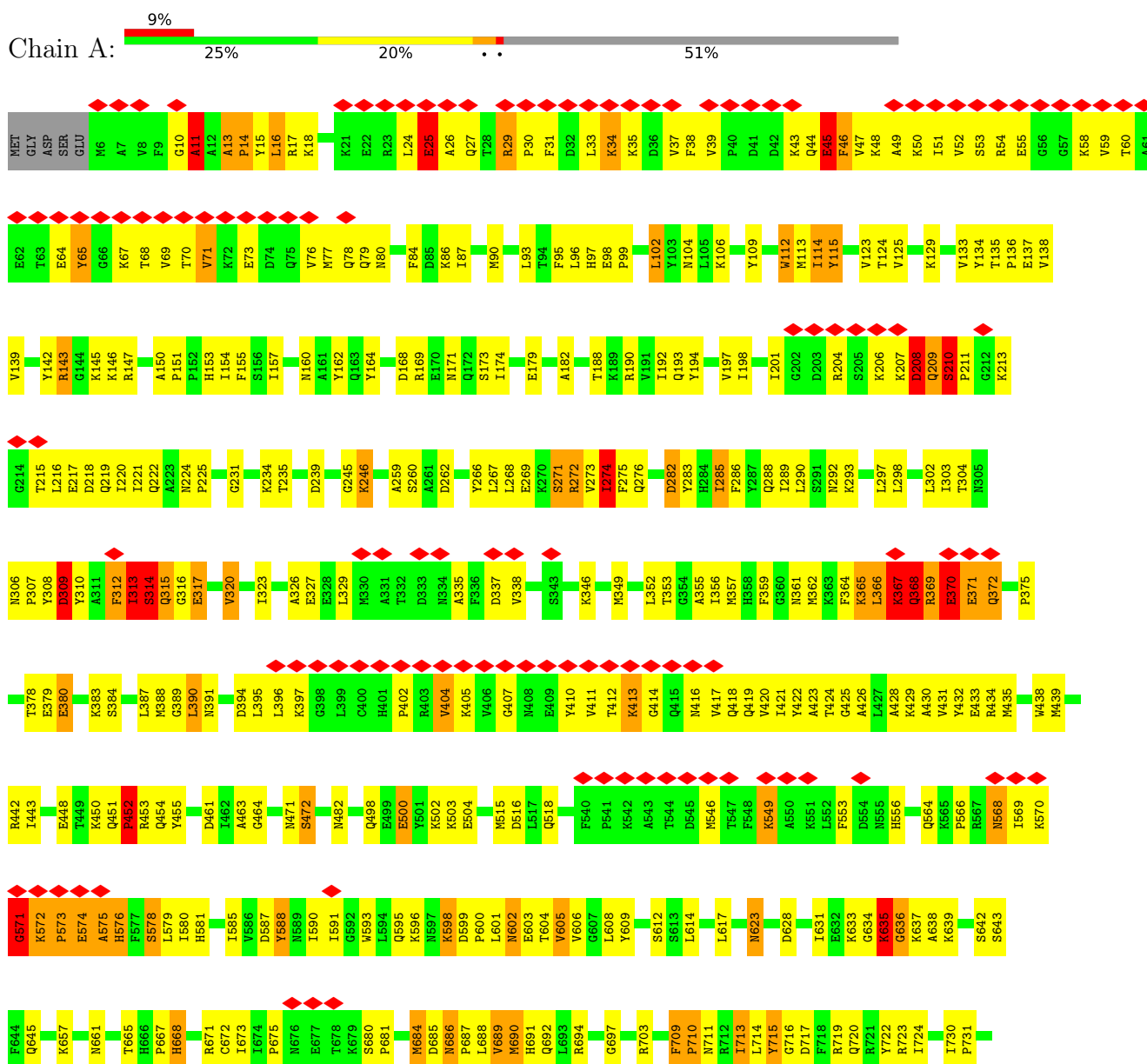
- Molecule 3 is a protein called Myosin regulatory light chain 2, ventricular/cardiac muscle isoform.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	160	Total	C	N	O	S	0	0
			1278	808	212	252	6		
3	F	160	Total	C	N	O	S	0	0
			1278	808	212	252	6		

3 Residue-property plots

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.

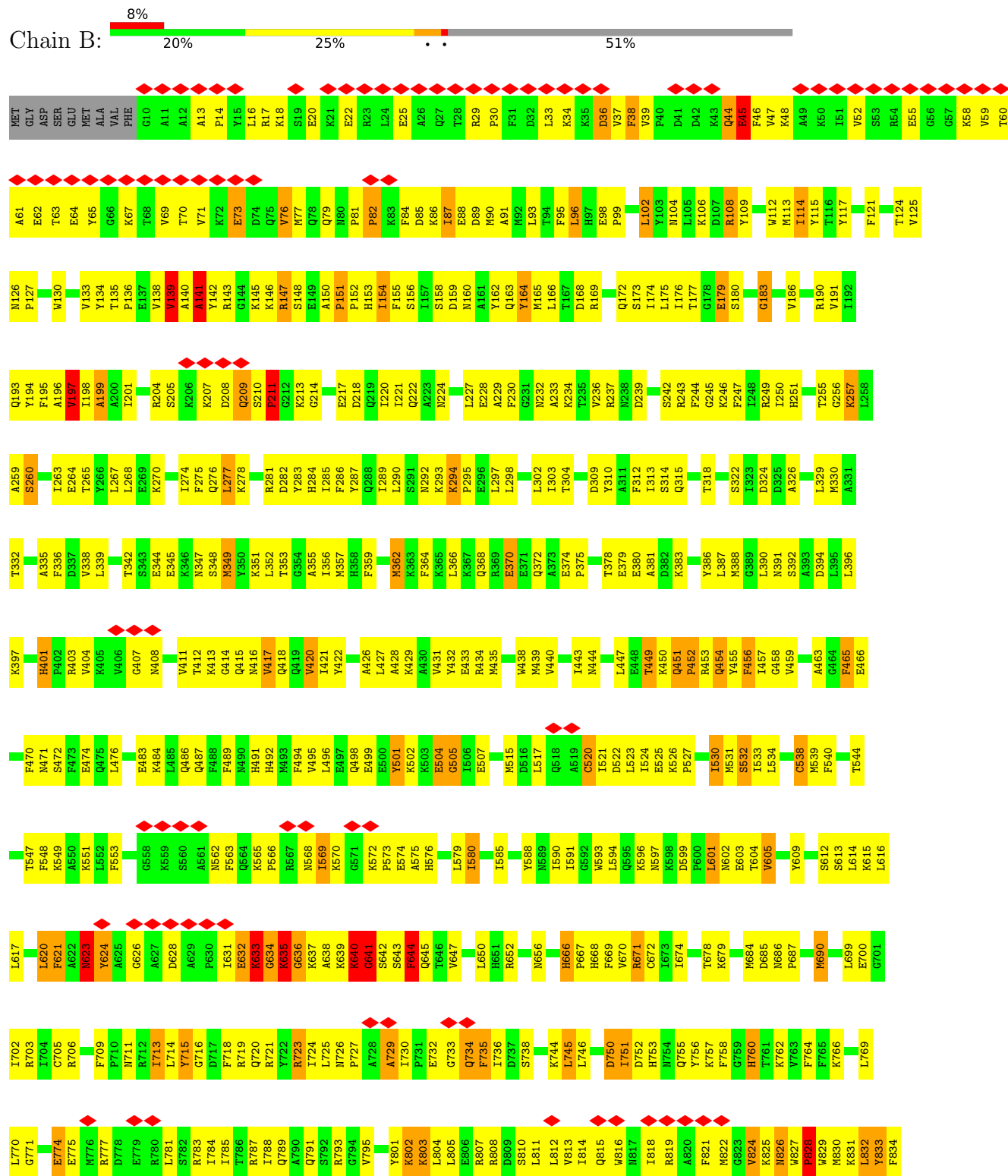
• Molecule 1: Myosin-7



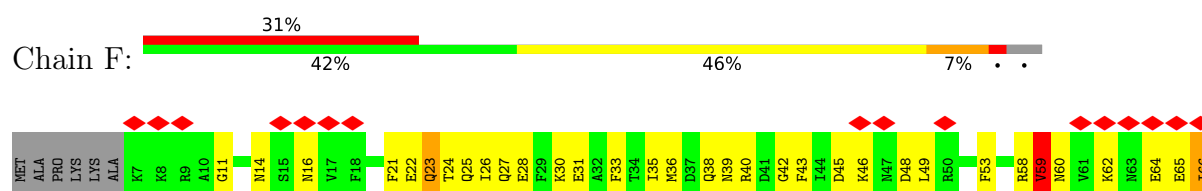


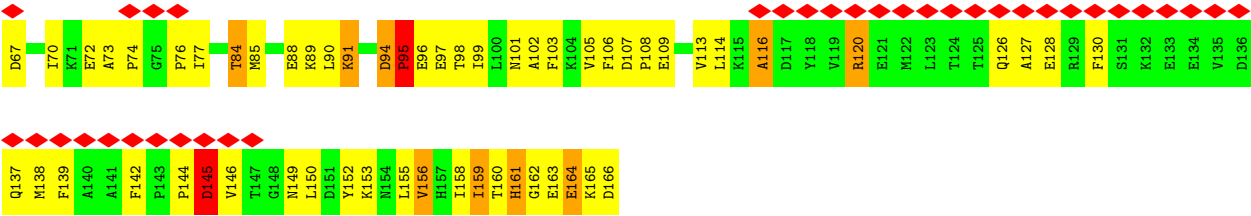
GLU	SER	GLN	VAL	ASN	LYS	LEU	ARG	ALA	LYS	SER	ARG	ASP	ILE	GLY	THR	LYS	GLY	LEU	ASN	GLU	GLU
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● Molecule 1: Myosin-7



- Molecule 2: Myosin light chain 3





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	10700	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI/PHILIPS CM120T	Depositor
Voltage (kV)	120	Depositor
Electron dose ($e^-/\text{\AA}^2$)	9	Depositor
Minimum defocus (nm)	1950	Depositor
Maximum defocus (nm)	1950	Depositor
Magnification	35000	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	12.488	Depositor
Minimum map value	-8.243	Depositor
Average map value	0.043	Depositor
Map value standard deviation	0.776	Depositor
Recommended contour level	1.5	Depositor
Map size ($\text{\AA}$)	950.4, 950.4, 950.4	wwPDB
Map dimensions	160, 160, 160	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	5.94, 5.94, 5.94	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	0.81	2/7851 (0.0%)	1.37	94/10556 (0.9%)
1	B	0.82	5/7819 (0.1%)	1.34	53/10513 (0.5%)
2	C	0.84	0/1231	1.32	3/1651 (0.2%)
2	D	1.07	6/1231 (0.5%)	1.38	7/1651 (0.4%)
3	E	1.04	5/1301 (0.4%)	1.37	8/1747 (0.5%)
3	F	0.84	0/1301	1.33	4/1747 (0.2%)
All	All	0.85	18/20734 (0.1%)	1.35	169/27865 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	20
1	B	0	19
2	C	0	2
3	E	0	3
3	F	0	4
All	All	0	48

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	129	ARG	CZ-NH1	-17.98	1.07	1.32
2	D	101	GLN	C-O	-15.26	1.02	1.24
1	B	666	HIS	CD2-NE2	-11.93	1.24	1.37
3	E	129	ARG	NE-CZ	11.07	1.45	1.33
2	D	105	ASN	C-N	-9.40	1.20	1.33

The worst 5 of 169 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	641	GLY	CA-C-O	-13.94	96.31	120.57
1	B	666	HIS	CE1-NE2-CD2	13.67	122.67	109.00
3	E	129	ARG	NE-CZ-NH2	-13.00	107.50	119.20
2	D	101	GLN	CA-C-O	-12.72	104.56	119.95
1	A	366	LEU	CA-C-N	12.01	143.32	121.70

There are no chirality outliers.

5 of 48 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	TYR	Sidechain
1	A	15	TYR	Sidechain
1	A	162	TYR	Sidechain
1	A	25	GLU	Mainchain
1	A	45	GLU	Mainchain

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	7704	0	7735	705	0
1	B	7673	0	7706	1090	0
2	C	1212	0	1182	125	0
2	D	1212	0	1184	236	0
3	E	1278	0	1241	180	0
3	F	1278	0	1239	238	0
All	All	20357	0	20287	2085	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 51.

The worst 5 of 2085 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:A:822:MET:HE1	3:E:138:MET:CE	1.27	1.64
1:B:804:LEU:CD2	2:D:63:ARG:HD3	1.21	1.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:714:LEU:HD21	1:B:396:LEU:CD1	1.15	1.55
1:B:831:LYS:HD3	3:F:163:GLU:CG	1.34	1.55
3:E:8:LYS:CD	3:F:59:VAL:HA	1.33	1.54

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	952/1935 (49%)	759 (80%)	141 (15%)	52 (6%)	1	15
1	B	948/1935 (49%)	781 (82%)	111 (12%)	56 (6%)	1	13
2	C	150/195 (77%)	132 (88%)	12 (8%)	6 (4%)	2	18
2	D	150/195 (77%)	133 (89%)	12 (8%)	5 (3%)	3	21
3	E	158/166 (95%)	111 (70%)	32 (20%)	15 (10%)	0	8
3	F	158/166 (95%)	115 (73%)	27 (17%)	16 (10%)	0	7
All	All	2516/4592 (55%)	2031 (81%)	335 (13%)	150 (6%)	2	13

5 of 150 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	ALA
1	A	27	GLN
1	A	208	ASP
1	A	272	ARG
1	A	313	ILE

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	831/1695 (49%)	813 (98%)	18 (2%)	45	64
1	B	828/1695 (49%)	811 (98%)	17 (2%)	47	65
2	C	133/167 (80%)	132 (99%)	1 (1%)	73	80
2	D	133/167 (80%)	131 (98%)	2 (2%)	57	72
3	E	137/141 (97%)	136 (99%)	1 (1%)	76	81
3	F	137/141 (97%)	133 (97%)	4 (3%)	37	58
All	All	2199/4006 (55%)	2156 (98%)	43 (2%)	48	66

5 of 43 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	770	LEU
2	C	186	GLU
1	B	775	GLU
1	B	837	LYS
2	D	185	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 62 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	415	GLN
3	E	25	GLN
1	B	691	HIS
2	D	191	HIS
3	F	149	ASN

5.3.3 RNA ⓘ

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.

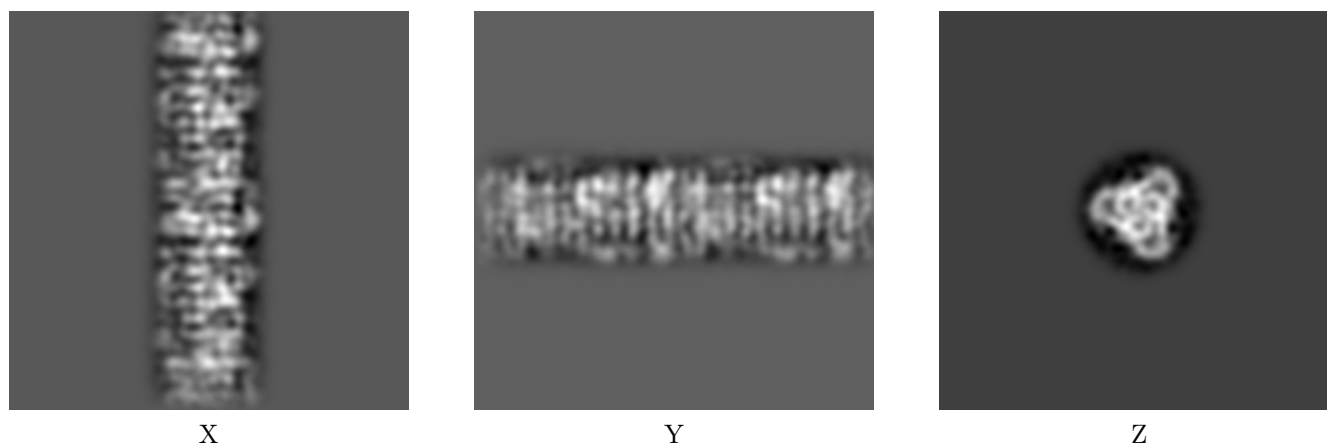
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-2240. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

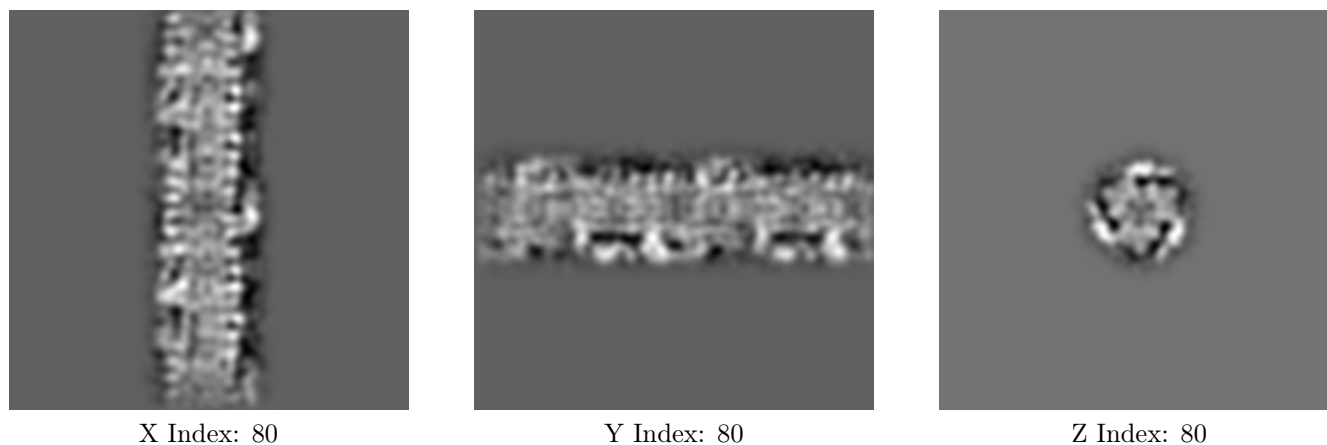
6.1.1 Primary map



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

6.2 Central slices [i](#)

6.2.1 Primary map



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

6.3 Largest variance slices [i](#)

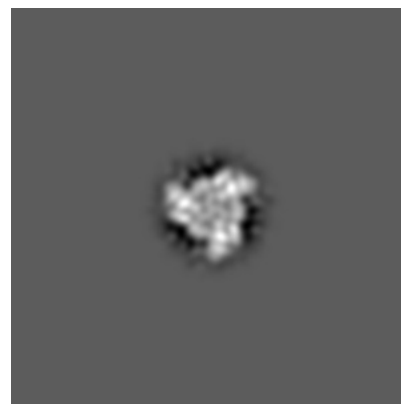
6.3.1 Primary map



X Index: 89



Y Index: 71

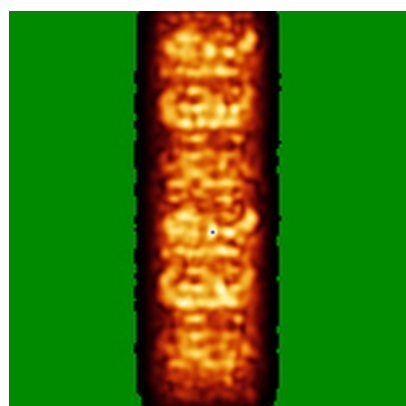


Z Index: 43

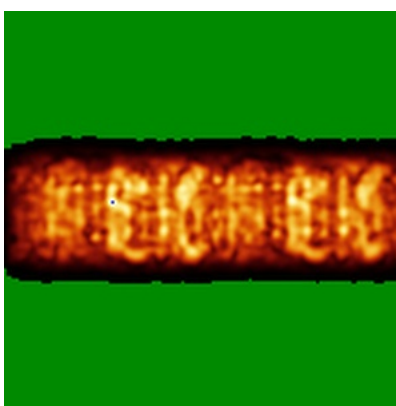
The images above show the largest variance slices of the map in three orthogonal directions.

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

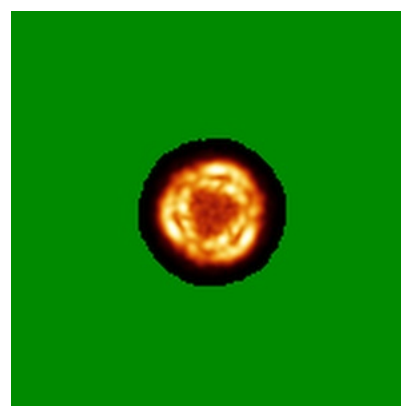
6.4.1 Primary map



X



Y

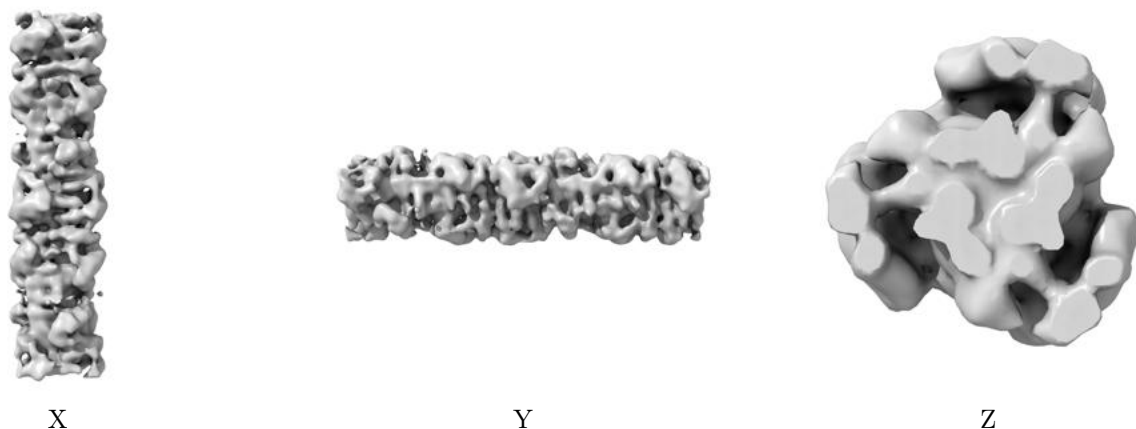


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

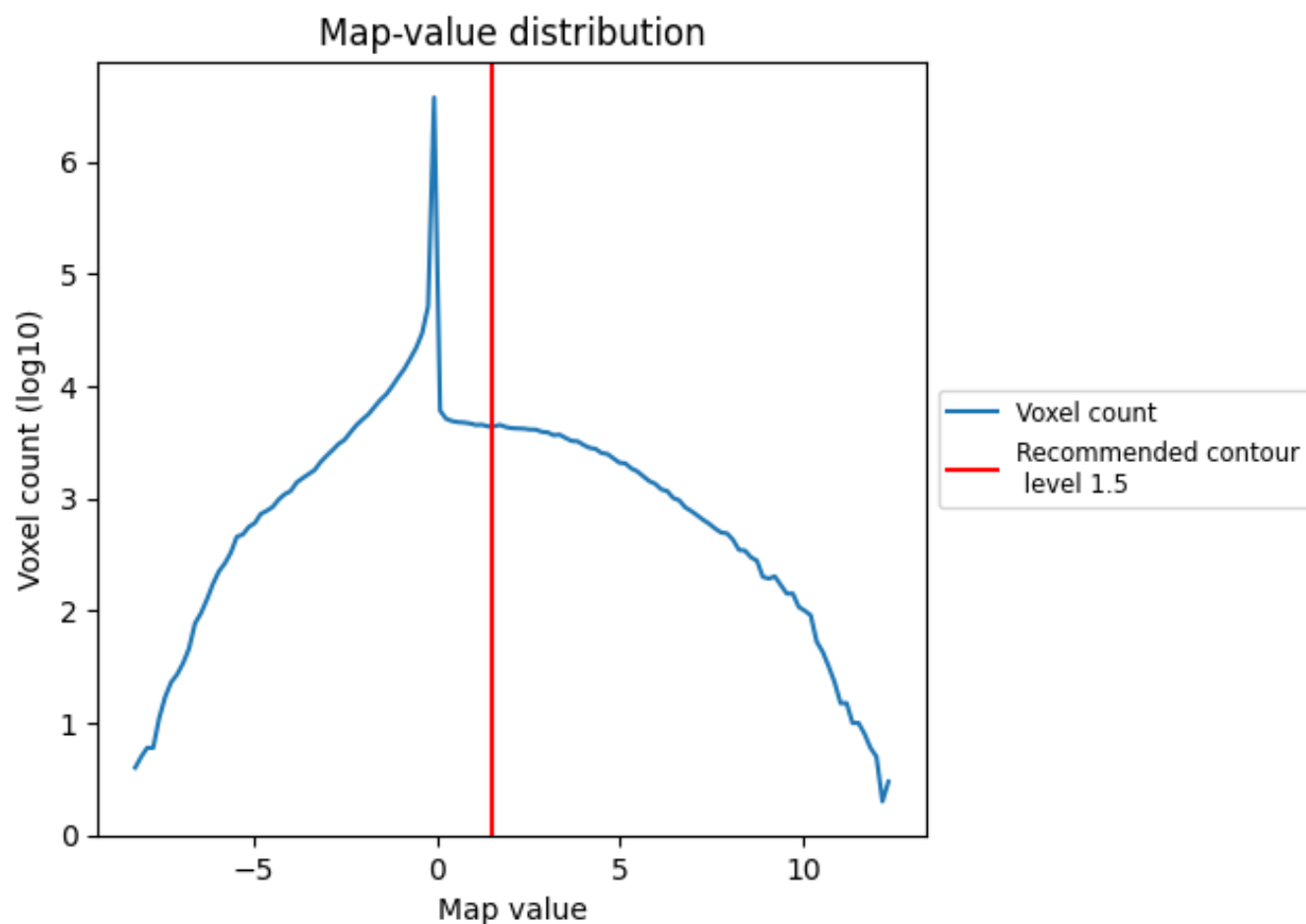
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

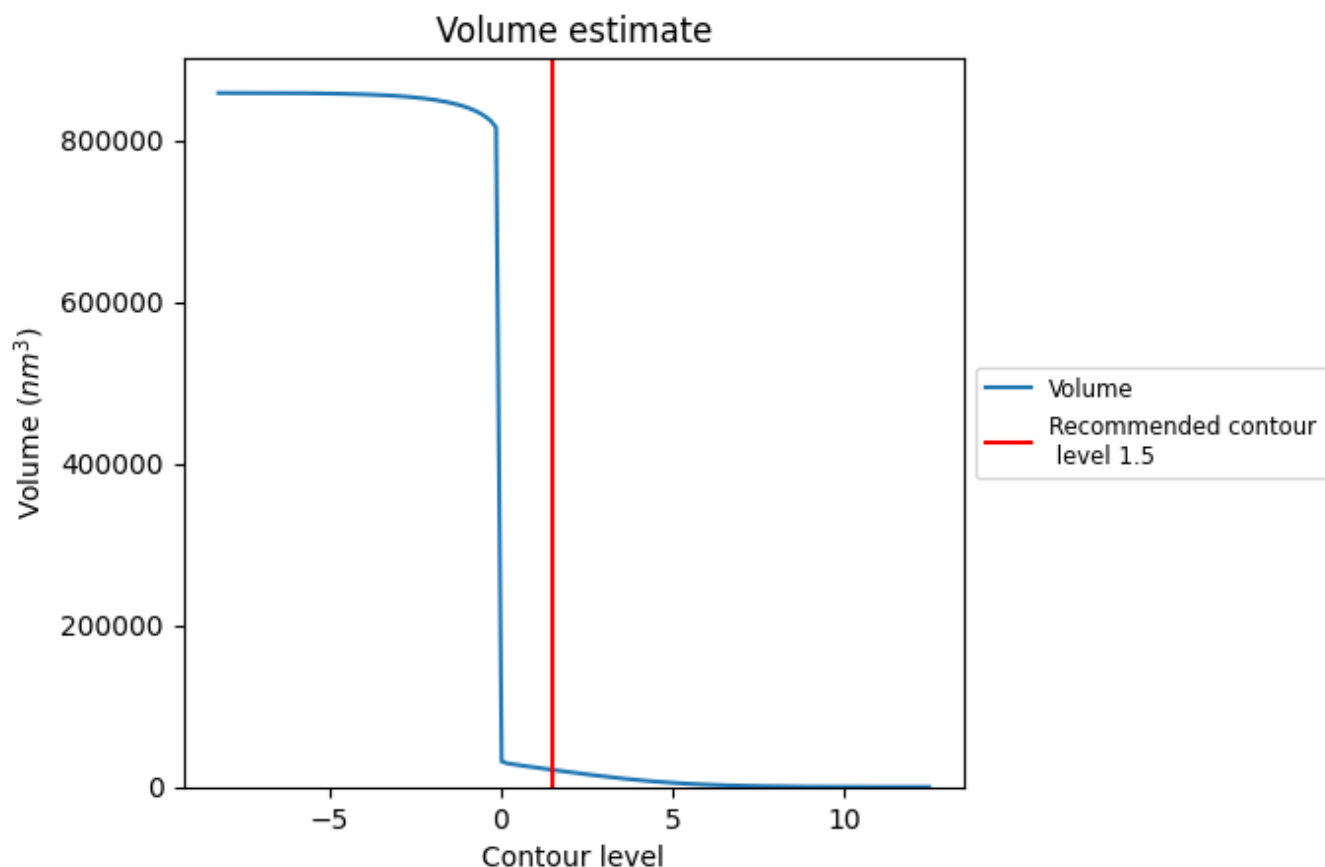
This section contains the results of statistical analysis of the map.

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.

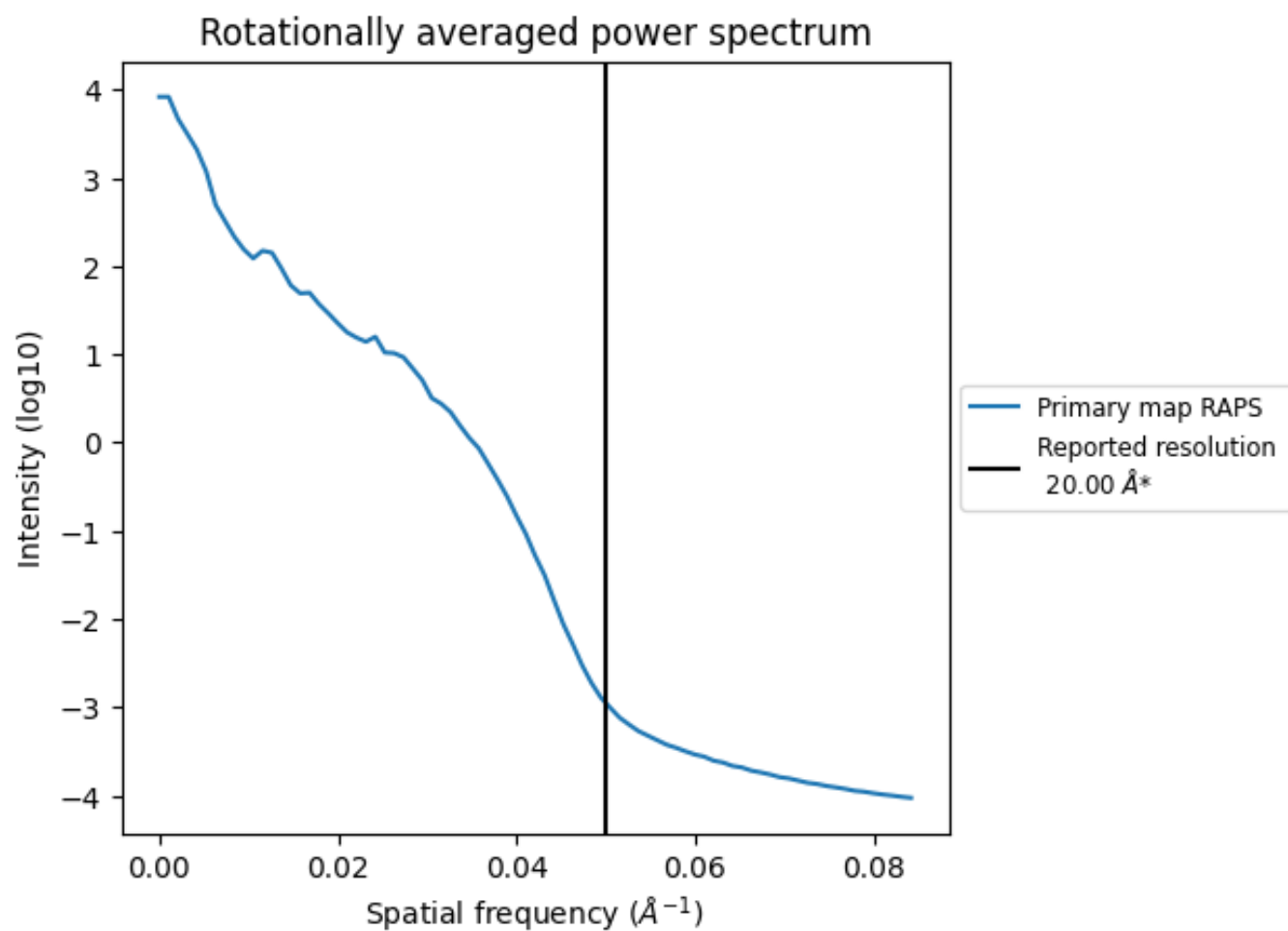
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 21179 nm^3 ; this corresponds to an approximate mass of 19131 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 0.050 $\text{\AA}^{-1}$

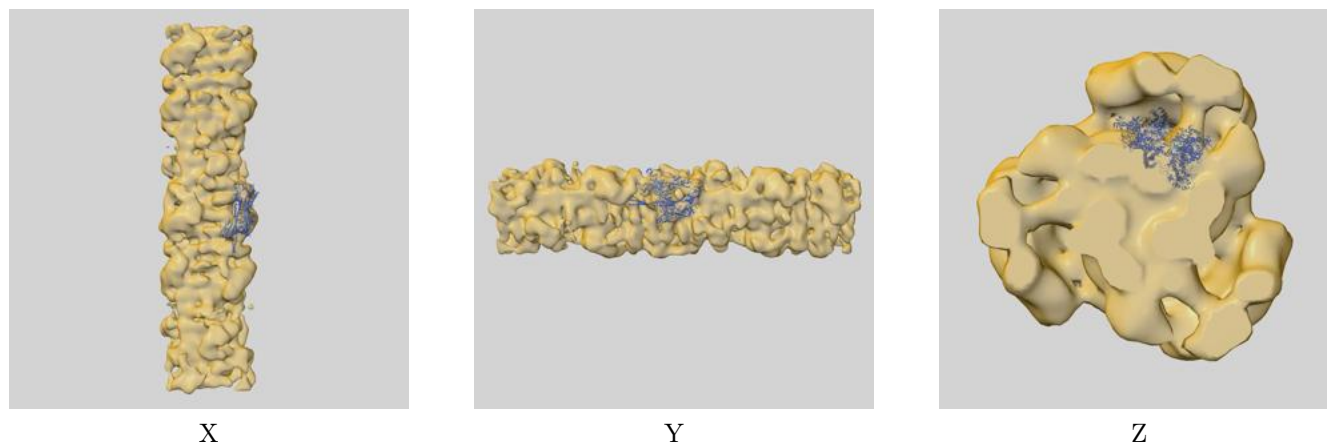
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

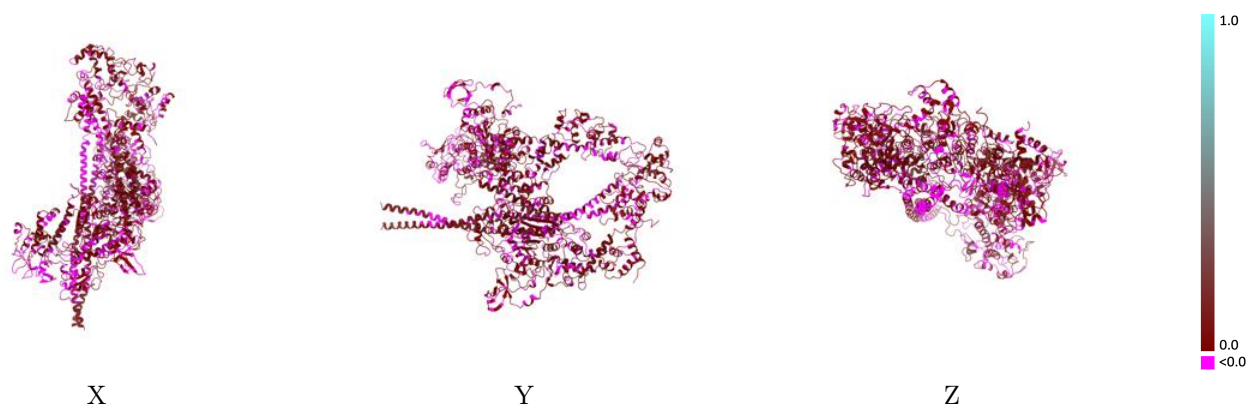
This section contains information regarding the fit between EMDB map EMD-2240 and PDB model 5TBY. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



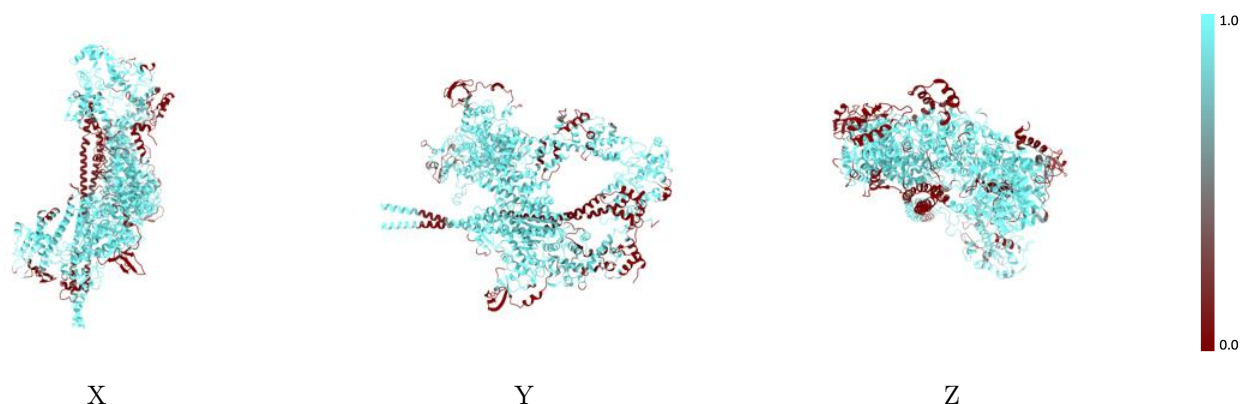
The images above show the 3D surface view of the map at the recommended contour level 1.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



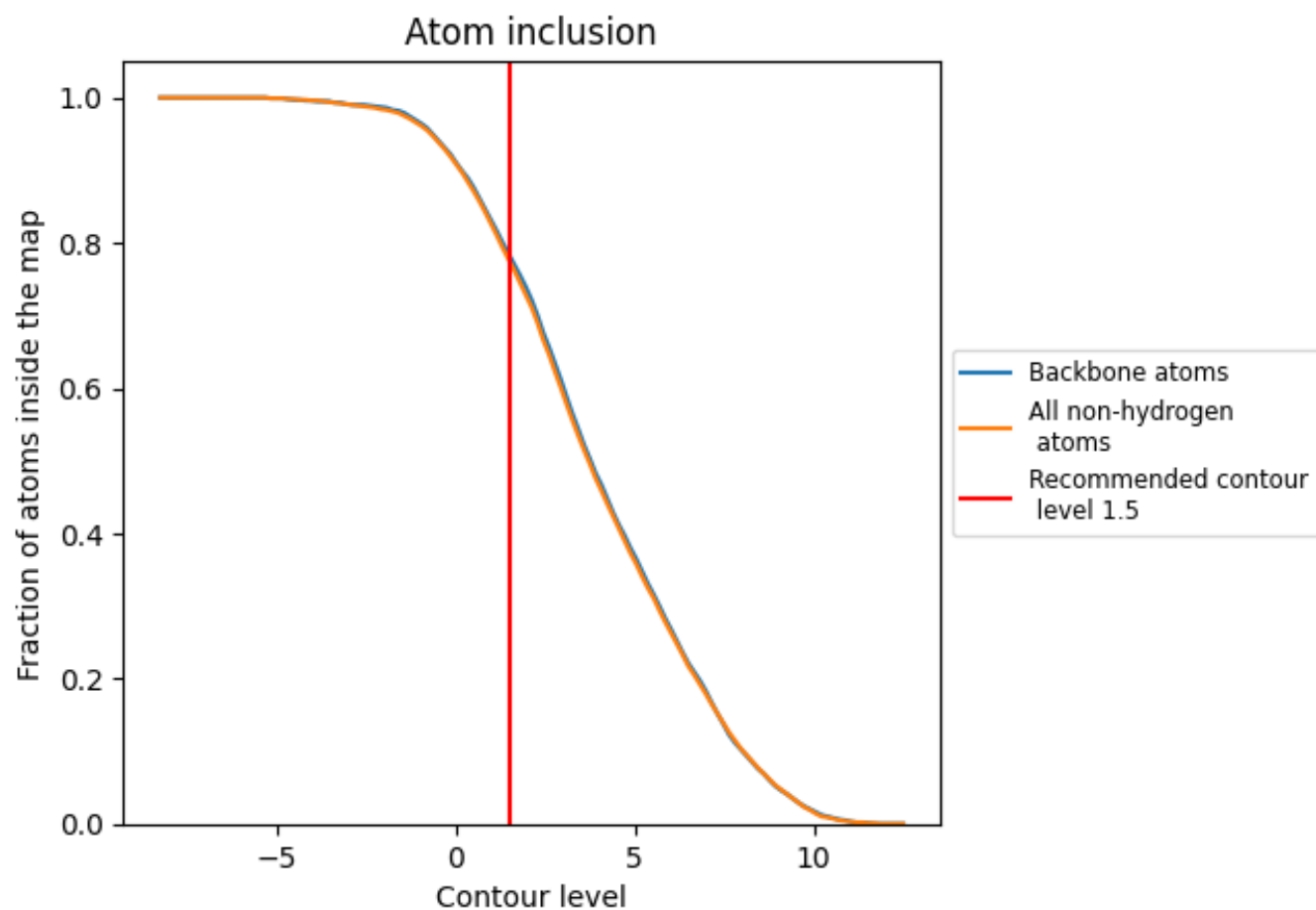
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 78% 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 (1.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7750	0.0320
A	0.8090	0.0230
B	0.8260	0.0380
C	0.5510	0.0350
D	0.4360	0.0430
E	0.9130	0.0340
F	0.6610	0.0420

1.0
0.0
-0.0