wwPDB EM Validation Summary Report

Nov 4, 2026 – 06:20 PM JST

PDB ID	:	43WO / pdb_000043wo
EMDB ID	:	EMD-82347
Title	:	Cryo-EM structure of the E. coli ArnA hexamer
Authors	:	Kim, S.; Yeo, H.K.; Lee, B.I.
Deposited on	:	2026-07-20
Resolution	:	3.24 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:

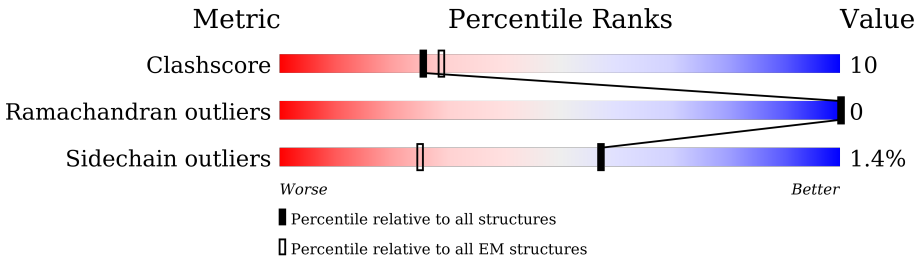
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

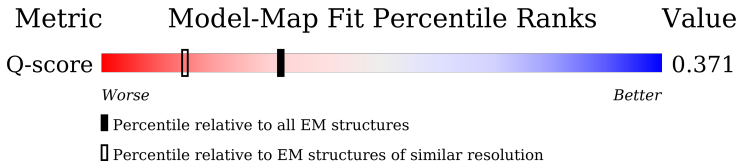
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.24 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	14594 (2.74 - 3.74)

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 <=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	653	70% 22% 8%
1	B	653	70% 23% 8%
1	C	653	68% 24% 8%
1	D	653	68% 24% 8%

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Mol	Chain	Length	Quality of chain
1	E	653	
1	F	653	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 28740 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 Bifunctional polymyxin resistance protein ArnA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	603	Total	C	N	O	S	0	0
			4789	3070	848	852	19		
1	B	603	Total	C	N	O	S	0	0
			4789	3070	848	852	19		
1	C	603	Total	C	N	O	S	0	0
			4789	3070	848	852	19		
1	D	603	Total	C	N	O	S	1	0
			4795	3074	850	852	19		
1	E	603	Total	C	N	O	S	0	0
			4789	3070	848	852	19		
1	F	603	Total	C	N	O	S	0	0
			4789	3070	848	852	19		

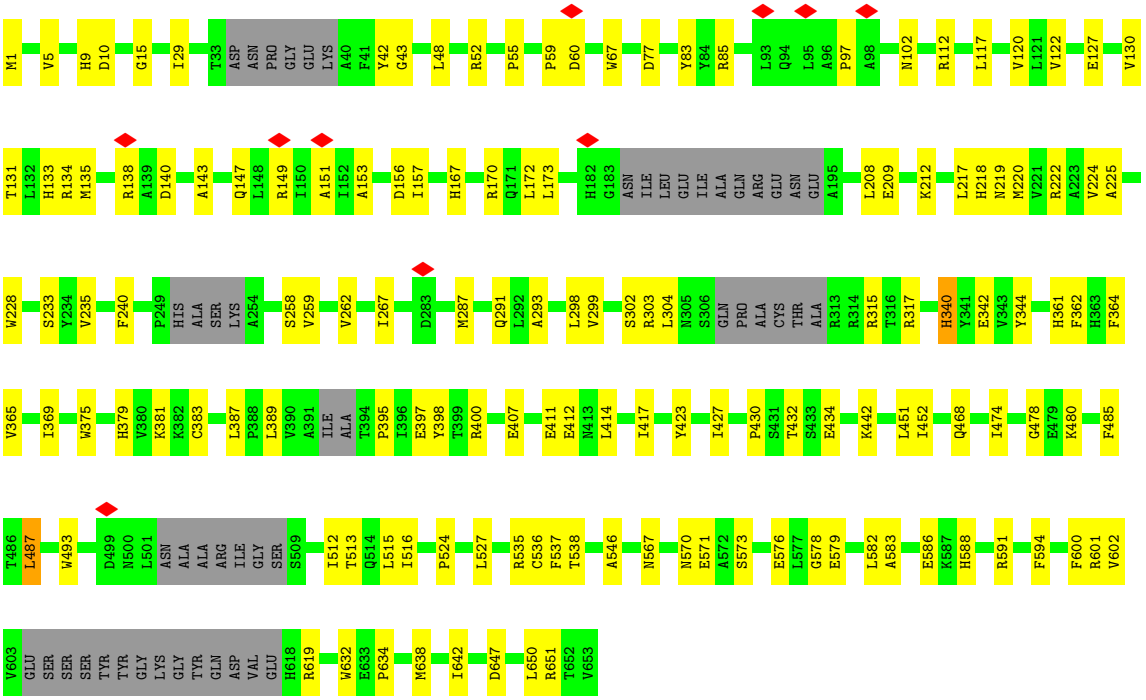
3 Residue-property plots (i)

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: Bifunctional polymyxin resistance protein ArnA

Chain A:

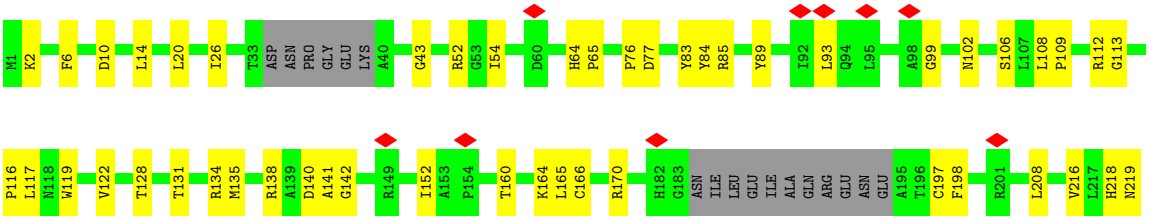


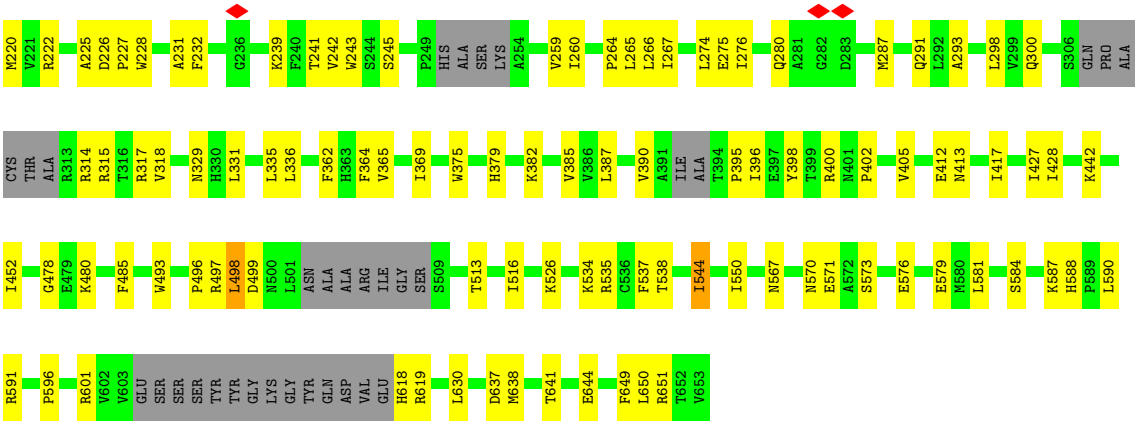


• Molecule 1: Bifunctional polymyxin resistance protein ArnA

Chain B:



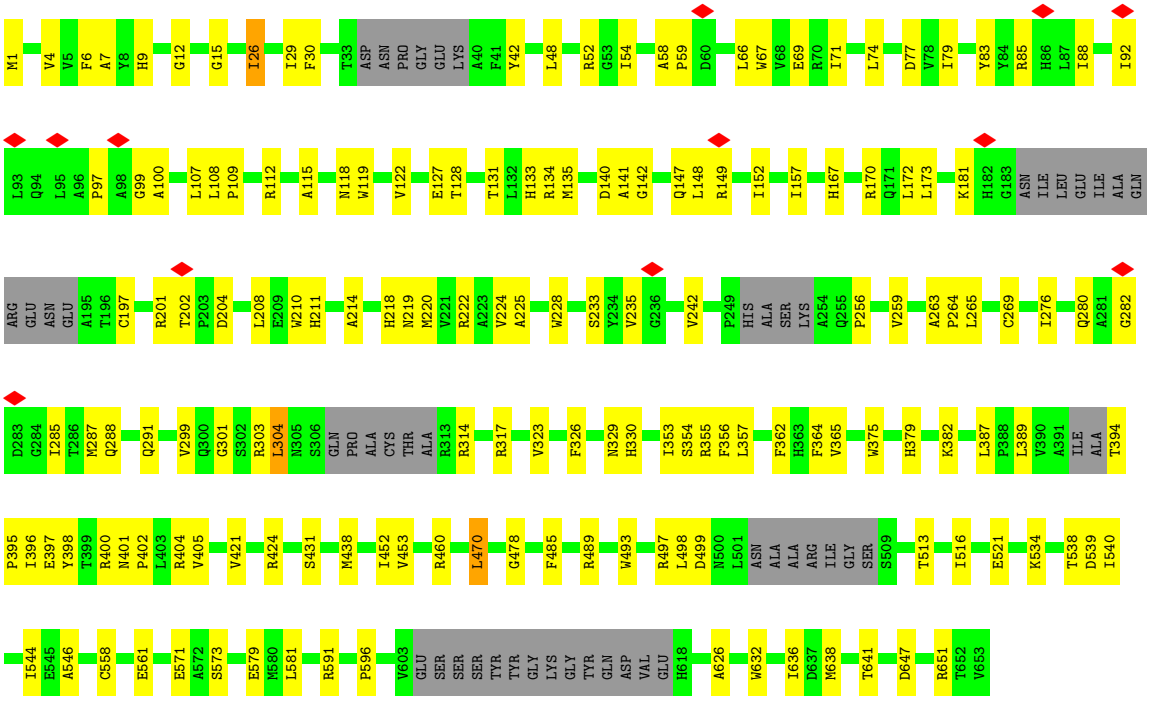




• Molecule 1: Bifunctional polymyxin resistance protein ArnA

Chain C:

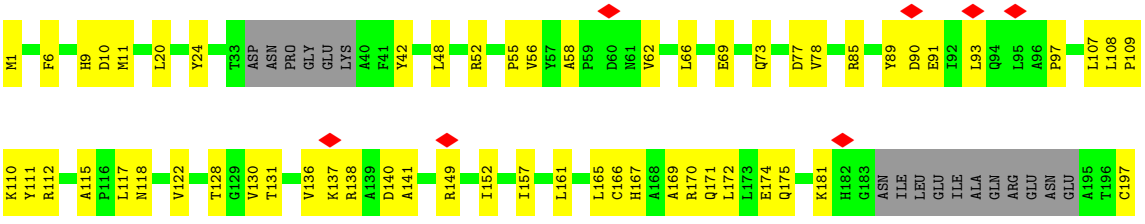


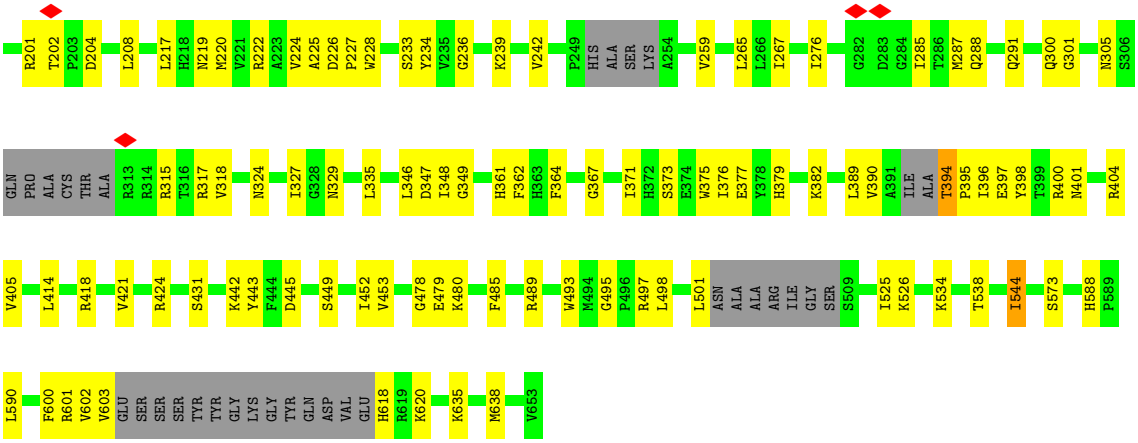


• Molecule 1: Bifunctional polymyxin resistance protein ArnA

Chain D:



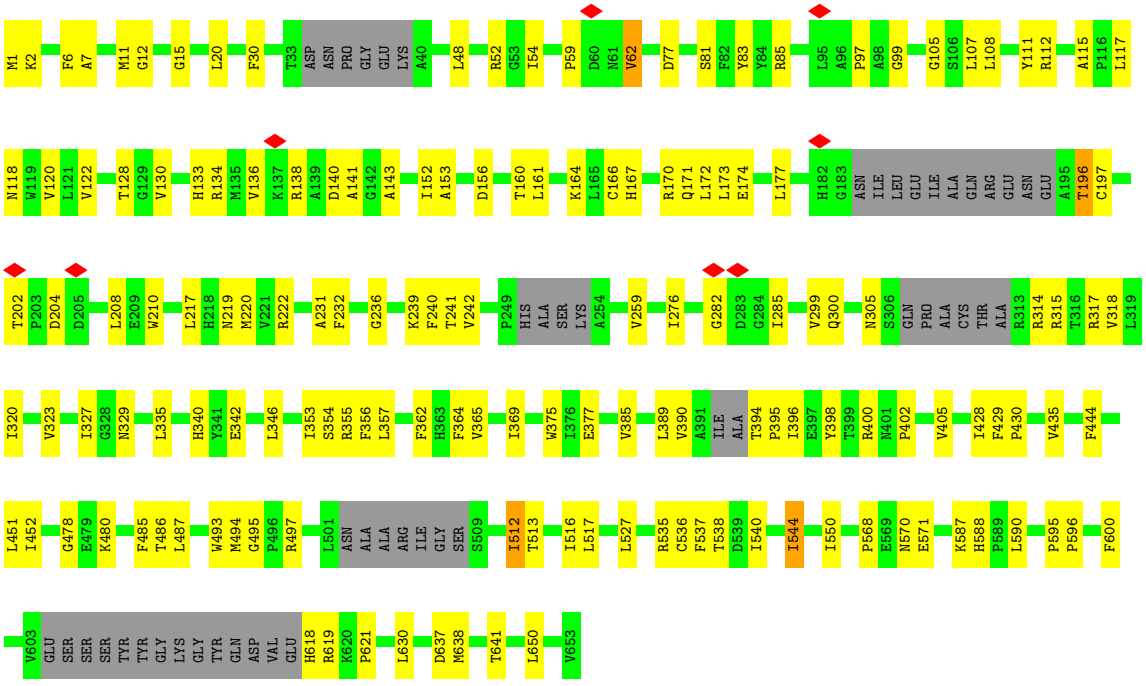




• Molecule 1: Bifunctional polymyxin resistance protein ArnA

Chain E:

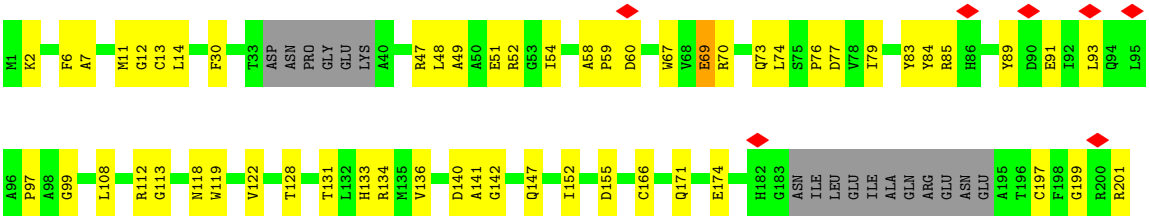


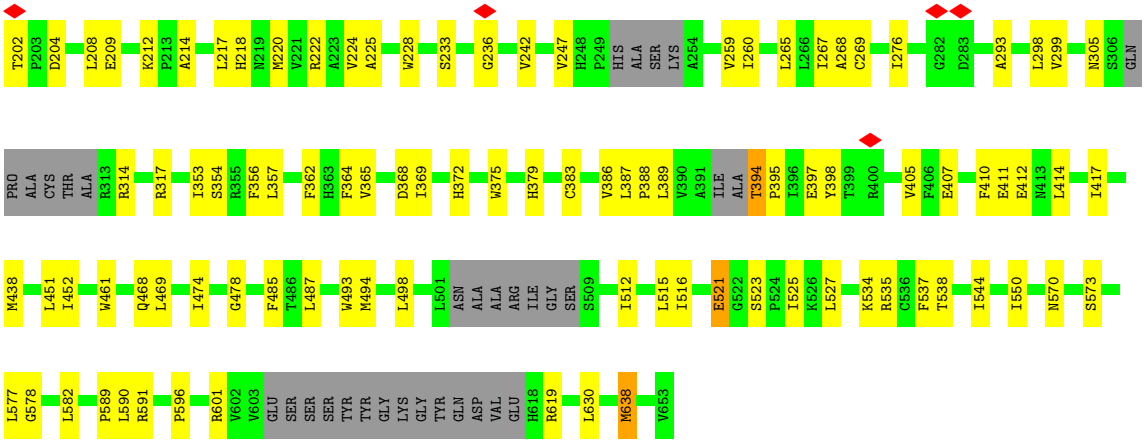


• Molecule 1: Bifunctional polymyxin resistance protein ArnA

Chain F:






T202 P203 D204 L208 E209 K212 P213 A214 L217 H218 W219 M220 V221 R222 A223 V224 A225 W228 S233 G236 V242 V247 H248 P249 HIS ALA SER LYS A254 V259 I260 L265 L266 I267 A268 C269 I276 G282 D283 A293 L298 V299 N305 S306 GLN
PRO ALA CYS THR ALA R313 R314 R317 I353 S354 R355 F356 L357 F362 H363 F364 V365 D368 I369 H372 W375 H379 C383 V386 L387 P388 L389 V390 A391 ILE ALA T394 P395 I396 E397 Y398 T399 R400 V405 F406 E407 F410 E411 E412 W413 L414 I417
M438 L451 I452 W461 Q468 L469 I474 G478 F485 T486 L487 W493 M494 L498 L501 ASN ALA ALA ARG ILE GLY SER S509 I512 L515 I516 E521 G522 S523 P524 I525 K526 L527 K534 R535 C536 F537 T538 I544 I550 N570 S573
L577 G578 L582 P589 L590 R591 P596 R601 V602 V603 GLU SER SER SER TYR TYR GLY LYS GLY TYR GLN ASP VAL GLU H618 R619 L630 M638 V653

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D3	Depositor
Number of particles used	260782	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1750	Depositor
Maximum defocus (nm)	2750	Depositor
Magnification	Not provided	
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.849	Depositor
Minimum map value	-0.455	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.024	Depositor
Recommended contour level	0.113	Depositor
Map size (Å)	314.496, 314.496, 314.496	wwPDB
Map dimensions	216, 216, 216	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.456, 1.456, 1.456	Depositor

5 Model quality

5.1 Standard geometry

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	0.15	0/4907	0.31	0/6658
1	B	0.16	0/4907	0.31	0/6658
1	C	0.15	0/4907	0.30	0/6658
1	D	0.16	0/4918	0.33	0/6673
1	E	0.15	0/4907	0.30	0/6658
1	F	0.15	0/4907	0.30	0/6658
All	All	0.15	0/29453	0.31	0/39963

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	4789	0	4736	93	0
1	B	4789	0	4736	109	0
1	C	4789	0	4734	100	0
1	D	4795	0	4739	107	0
1	E	4789	0	4736	102	0
1	F	4789	0	4736	99	0
All	All	28740	0	28417	588	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 10.

The worst 5 of 588 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:99:GLY:HA3	1:C:134:ARG:HH12	1.36	0.90
1:B:535:ARG:HH11	1:B:619:ARG:HG2	1.50	0.75
1:C:112:ARG:HG2	1:C:140:ASP:HB2	1.67	0.75
1:C:591:ARG:HH22	1:C:596:PRO:HA	1.52	0.75
1:C:122:VAL:HG23	1:C:220:MET:HB3	1.69	0.74

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	587/653 (90%)	553 (94%)	34 (6%)	0	100	100
1	B	587/653 (90%)	563 (96%)	24 (4%)	0	100	100
1	C	587/653 (90%)	560 (95%)	27 (5%)	0	100	100
1	D	588/653 (90%)	558 (95%)	30 (5%)	0	100	100
1	E	587/653 (90%)	548 (93%)	39 (7%)	0	100	100
1	F	587/653 (90%)	554 (94%)	33 (6%)	0	100	100
All	All	3523/3918 (90%)	3336 (95%)	187 (5%)	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	508/552 (92%)	503 (99%)	5 (1%)	68	78
1	B	508/552 (92%)	503 (99%)	5 (1%)	68	78
1	C	508/552 (92%)	501 (99%)	7 (1%)	59	74
1	D	509/552 (92%)	501 (98%)	8 (2%)	55	72
1	E	508/552 (92%)	497 (98%)	11 (2%)	45	68
1	F	508/552 (92%)	500 (98%)	8 (2%)	55	72
All	All	3049/3312 (92%)	3005 (99%)	44 (1%)	57	74

5 of 44 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	340	HIS
1	E	544	ILE
1	E	369	ILE
1	E	394	THR
1	F	133	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	94	GLN
1	E	162	HIS
1	E	133	HIS
1	E	171	GLN
1	C	171	GLN

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.

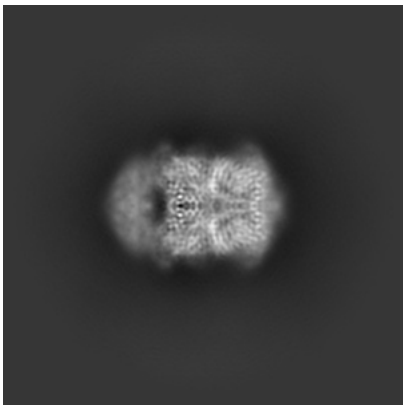
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-82347. These allow visual inspection of the internal detail of the map and identification of artifacts.

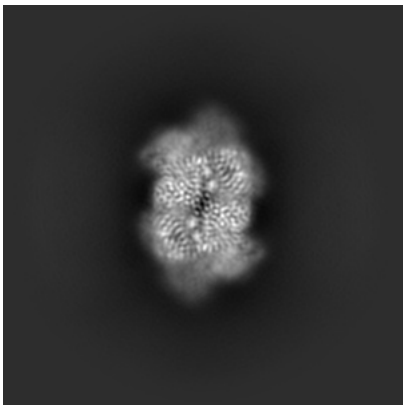
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

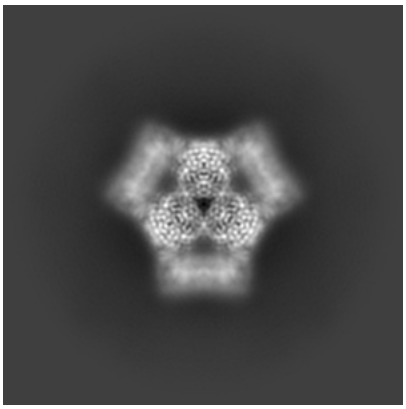
6.1.1 Primary map



X

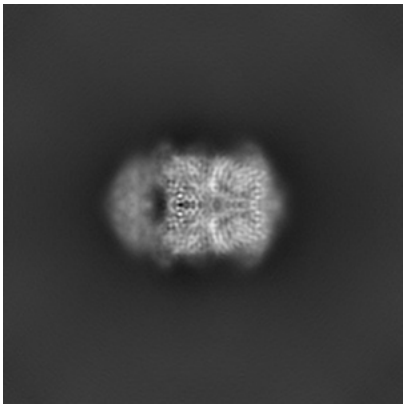


Y

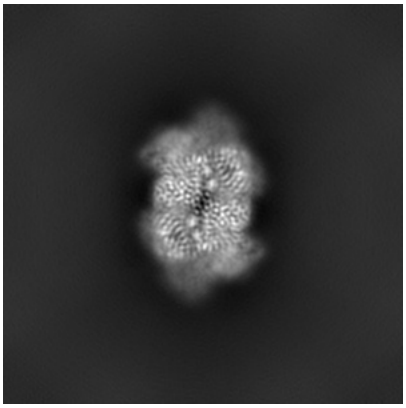


Z

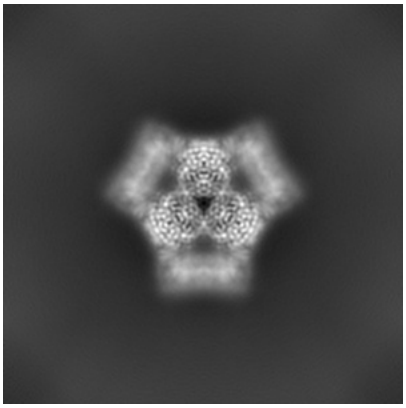
6.1.2 Raw map



X



Y

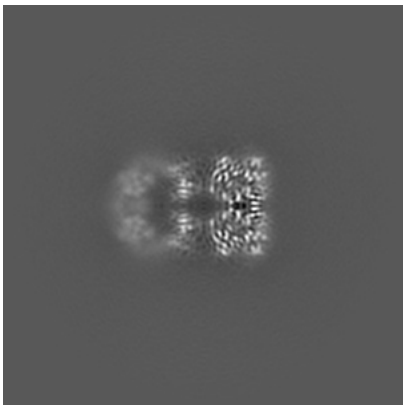


Z

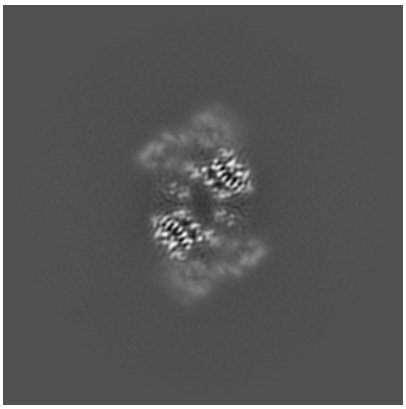
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

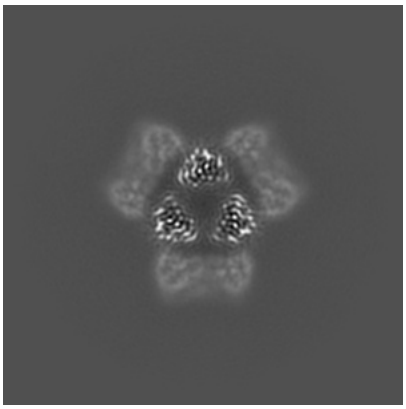
6.2.1 Primary map



X Index: 108

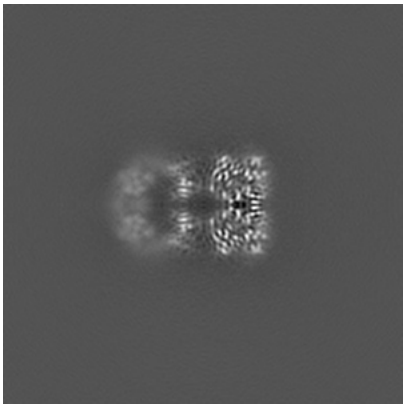


Y Index: 108

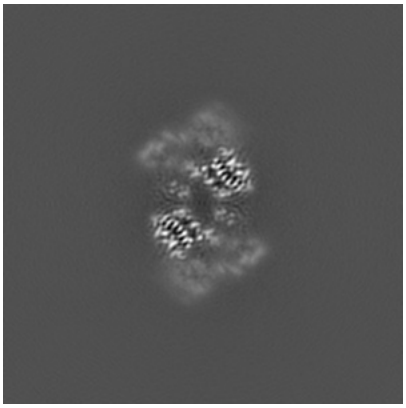


Z Index: 108

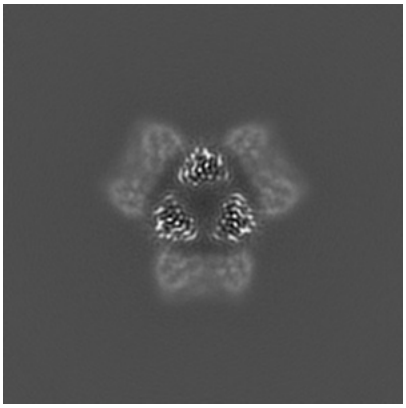
6.2.2 Raw map



X Index: 108



Y Index: 108

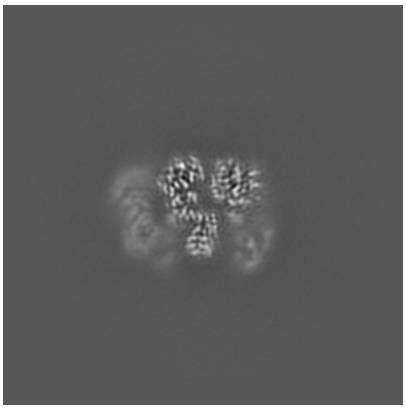


Z Index: 108

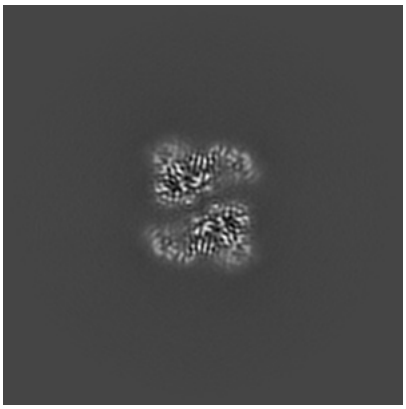
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

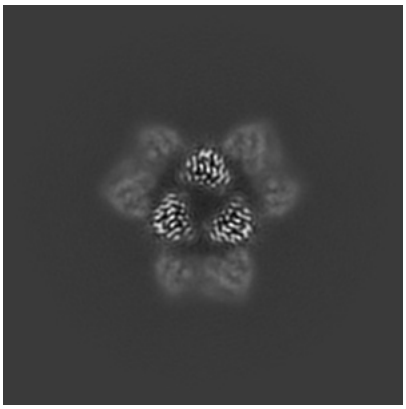
6.3.1 Primary map



X Index: 98

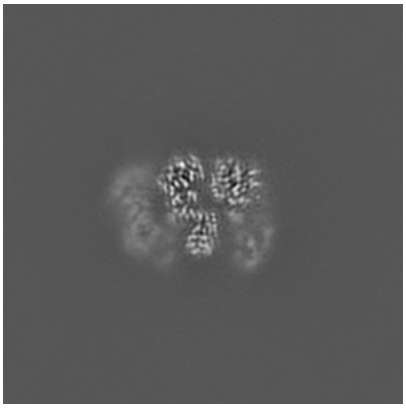


Y Index: 94

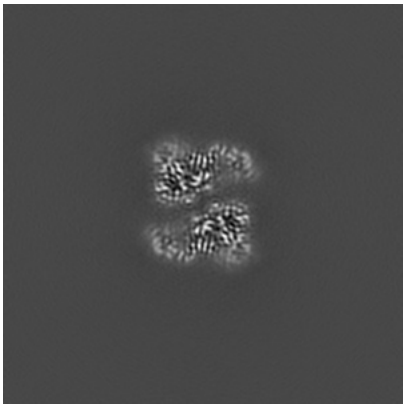


Z Index: 105

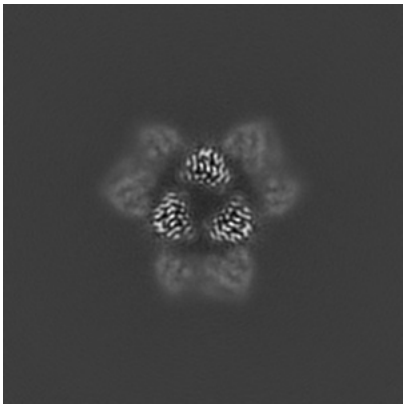
6.3.2 Raw map



X Index: 98



Y Index: 94

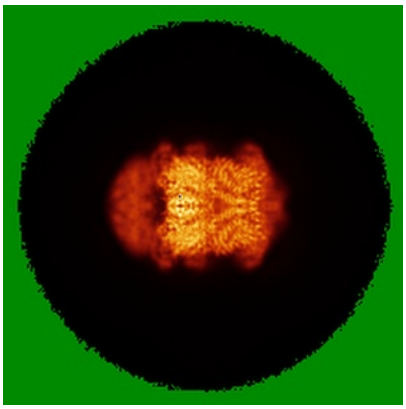


Z Index: 105

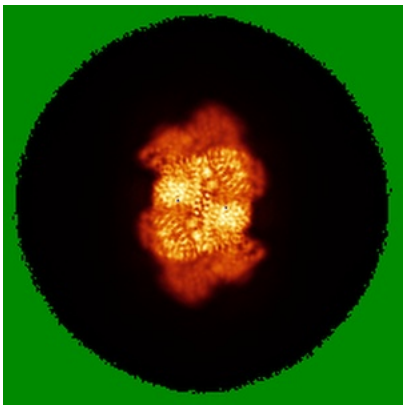
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

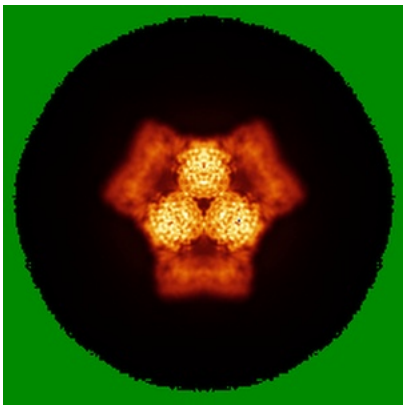
6.4.1 Primary map



X

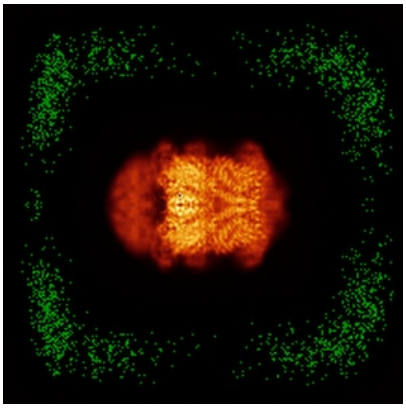


Y

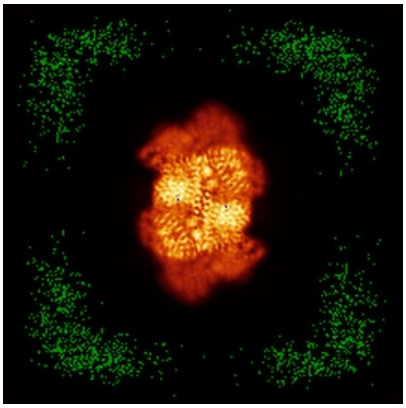


Z

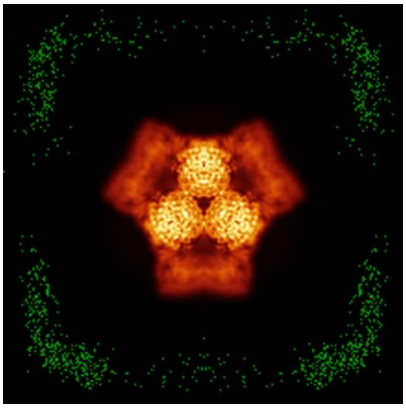
6.4.2 Raw map



X



Y

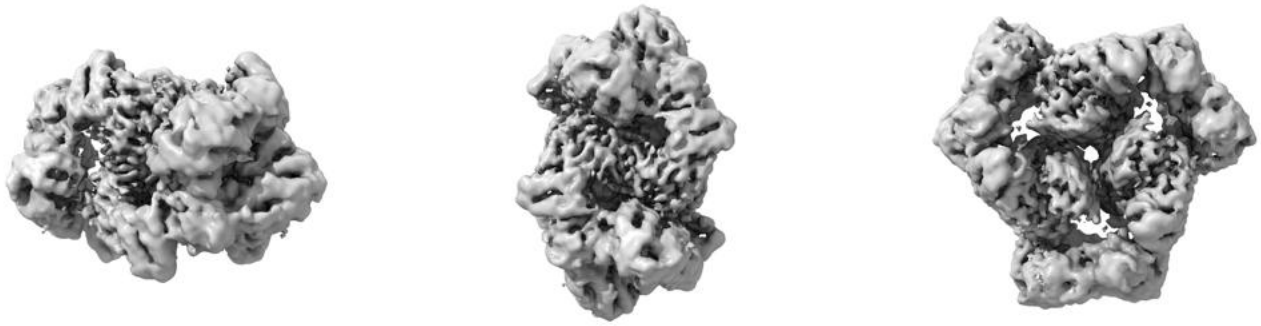


Z

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)

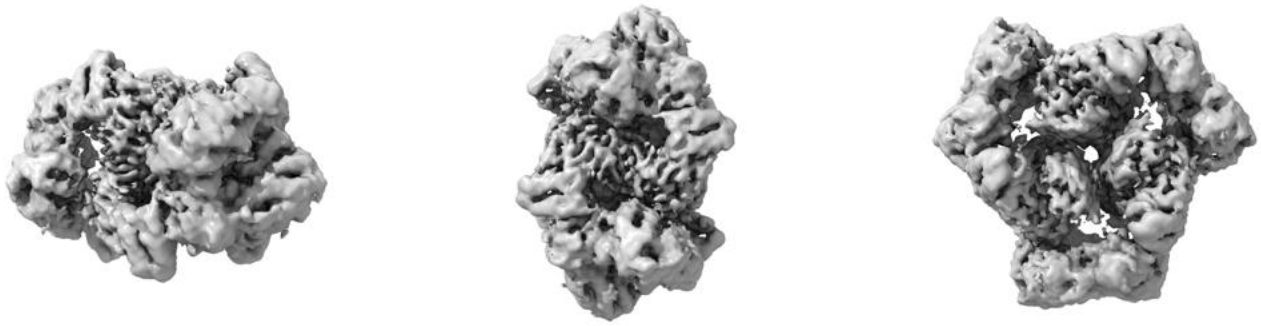
6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.113. 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.

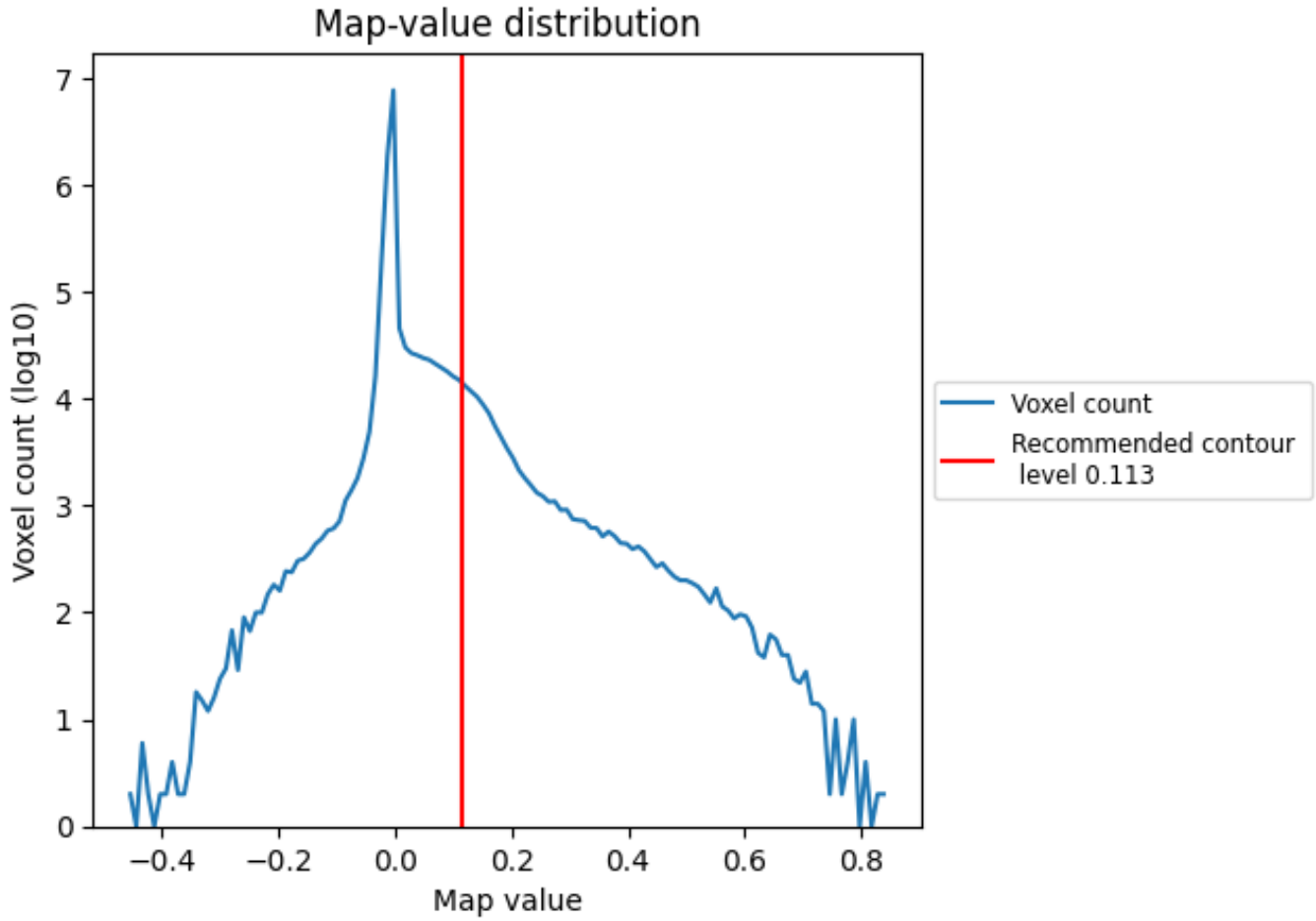
6.6 Mask visualisation (i)

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

7 Map analysis

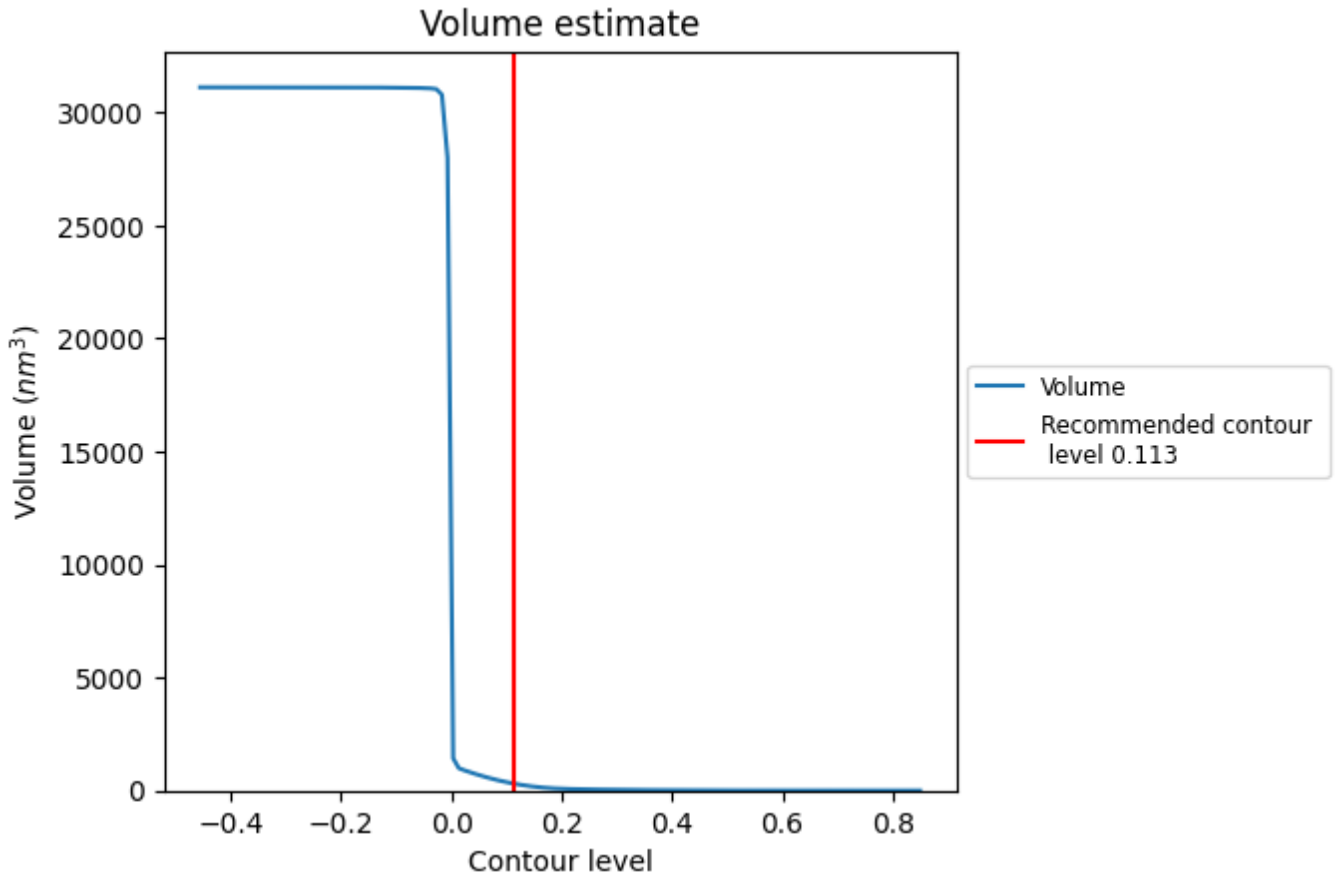
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

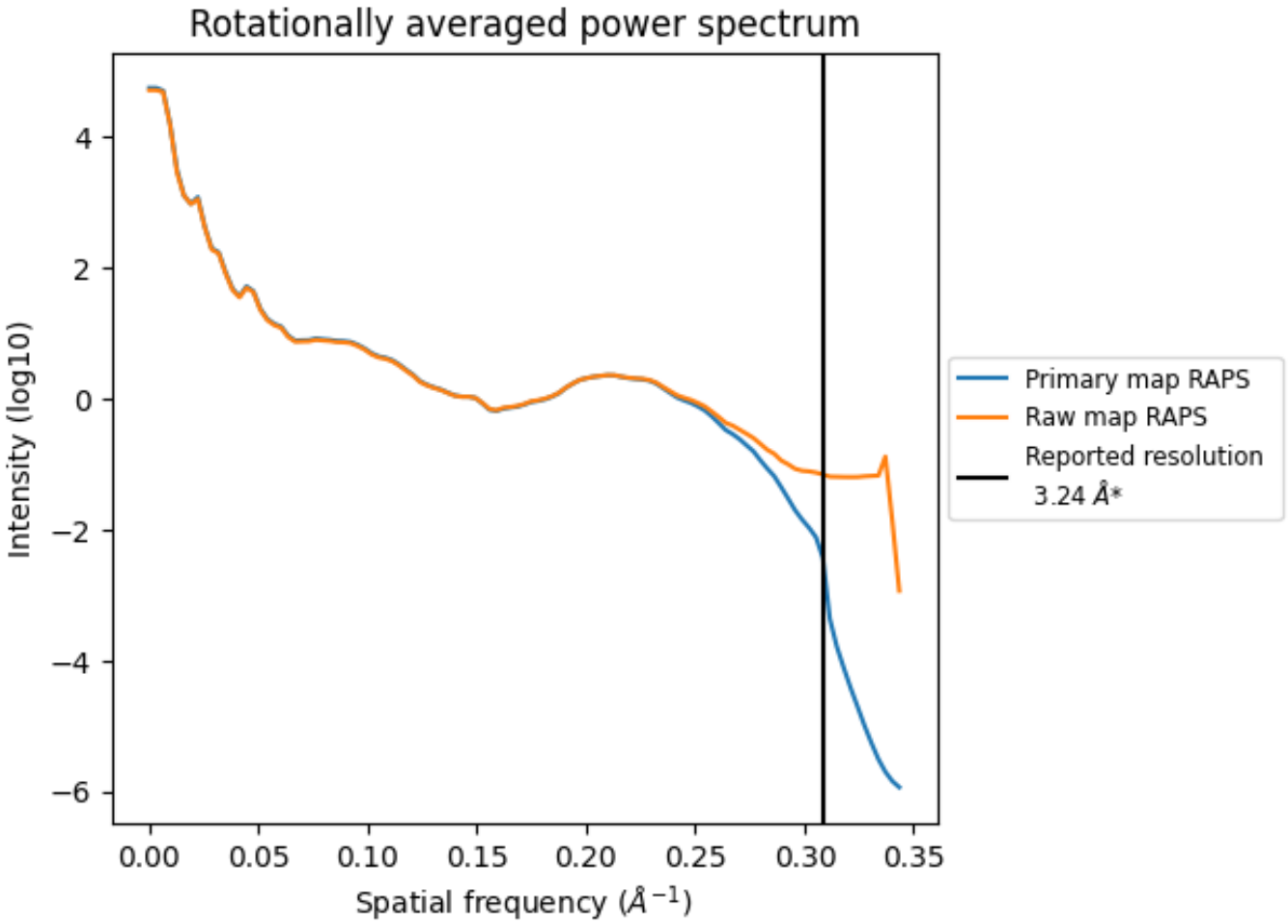
7.2 Volume estimate (i)



The volume at the recommended contour level is 310 nm^3; this corresponds to an approximate mass of 280 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)

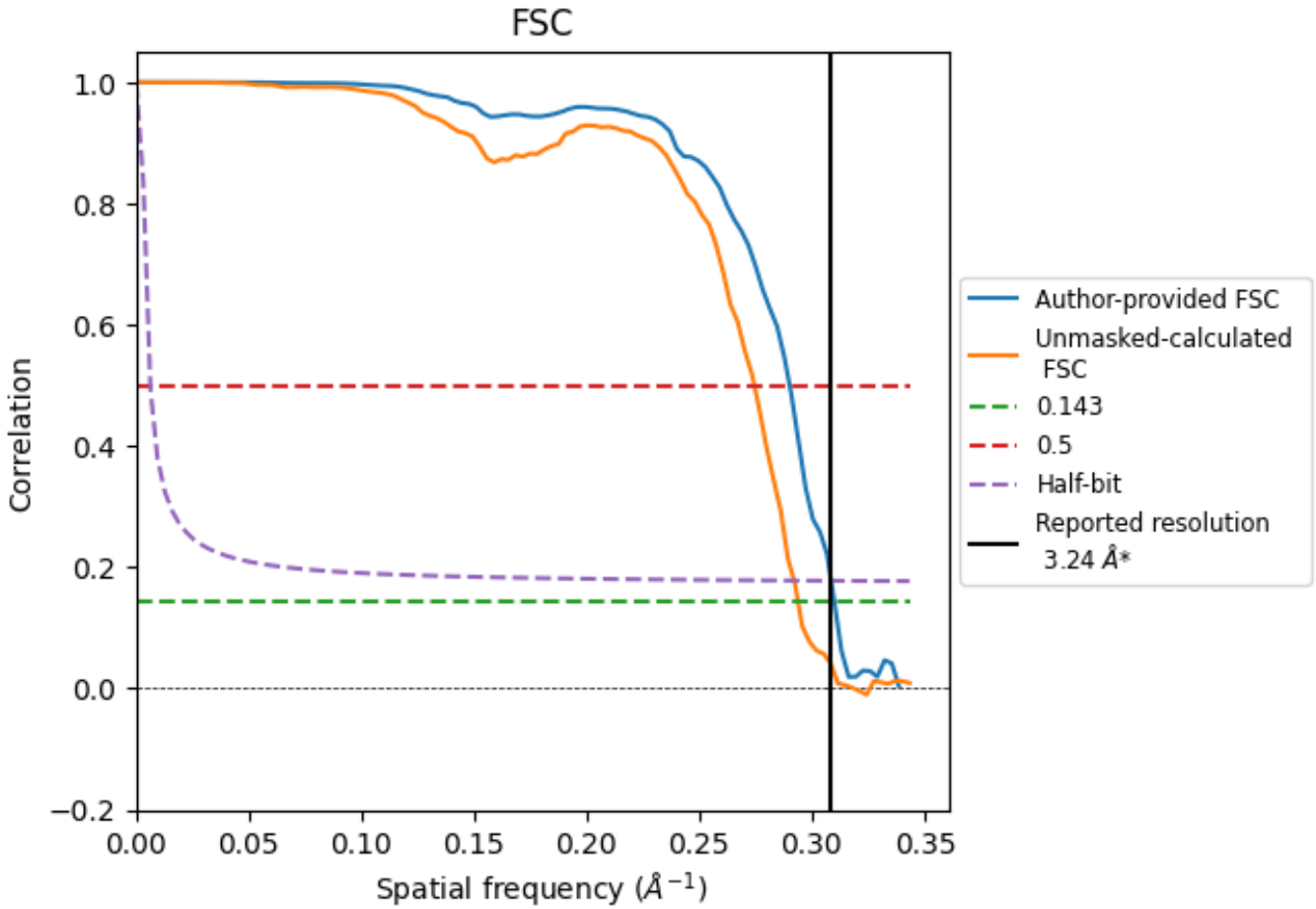


*Reported resolution corresponds to spatial frequency of 0.309 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.309 Å^{-1}

8.2 Resolution estimates (i)

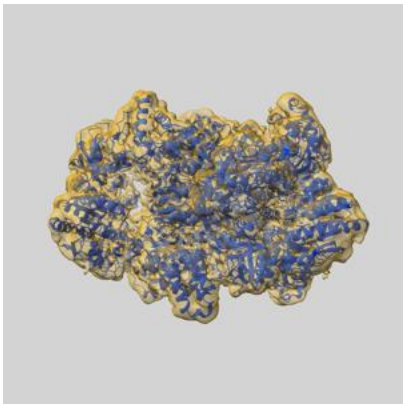
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.24	-	-
Author-provided FSC curve	3.23	3.45	3.24
Unmasked-calculated*	3.40	3.64	3.42

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

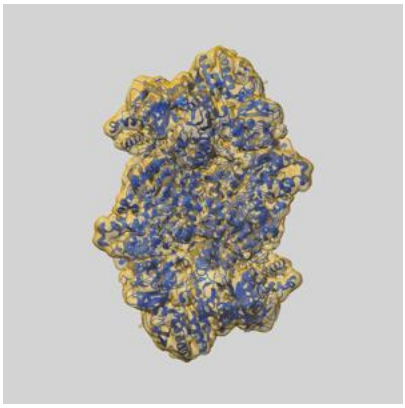
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-82347 and PDB model 43WO. Per-residue inclusion information can be found in section 3 on page 5.

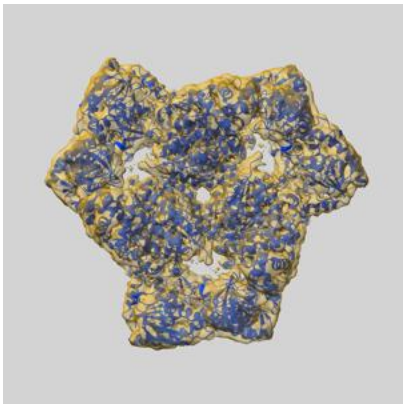
9.1 Map-model overlay



X



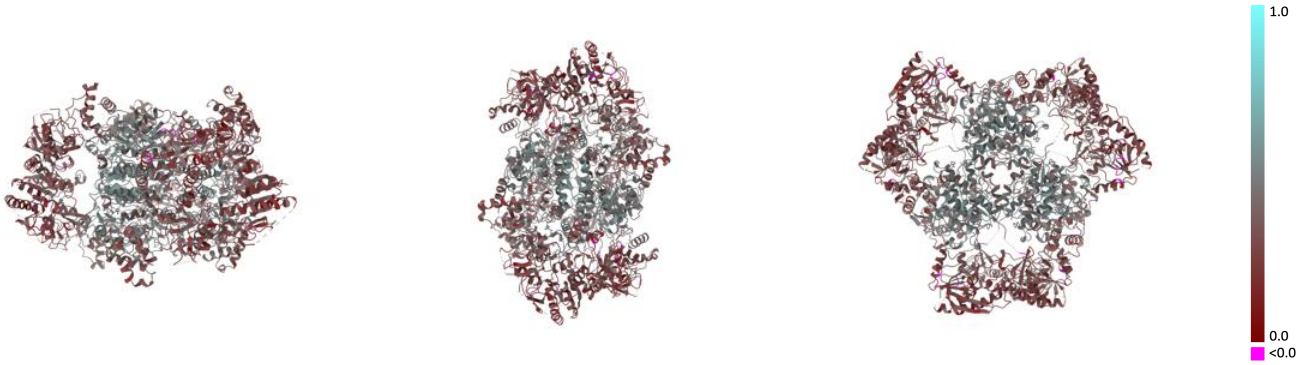
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.113 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.

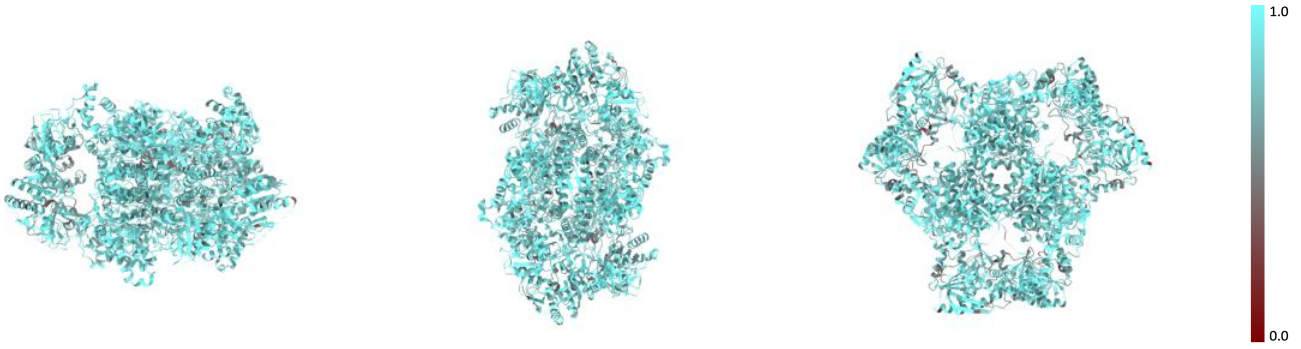
9.2 Q-score mapped to coordinate model



X Y Z

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.

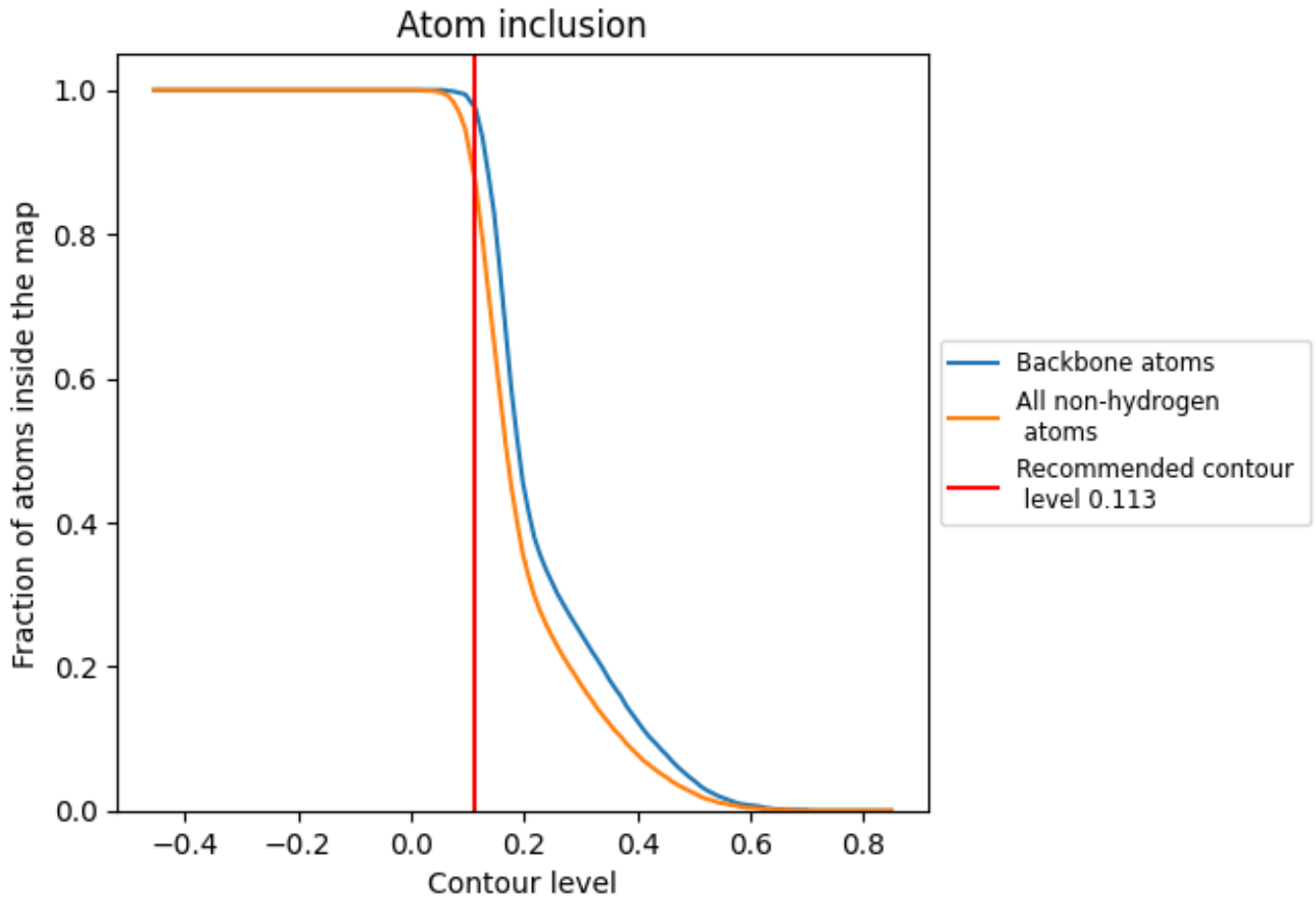
9.3 Atom inclusion mapped to coordinate model



X Y Z

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 (0.113).

9.4 Atom inclusion (i)



At the recommended contour level, 98% of all backbone atoms, 87% 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 (0.113) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8730	0.3710
A	0.8720	0.3700
B	0.8740	0.3740
C	0.8770	0.3720
D	0.8690	0.3700
E	0.8730	0.3680
F	0.8710	0.3700

