Full wwPDB EM Validation Report

Mar 8, 2026 – 12:49 PM UTC

PDB ID	:	9HX4 / pdb_00009hx4
EMDB ID	:	EMD-52457
Title	:	Amyloid fibril of apolipoprotein A-IV
Authors	:	Aibara, S.; Kassner, A.; Wong, E.; Klingel, K.; Papworth, M.; Althage, M.; Wang, Q.D.; Correia, C.; Milting, H.; Oliveira, T.M.
Deposited on	:	2025-01-06
Resolution	:	3.30 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:

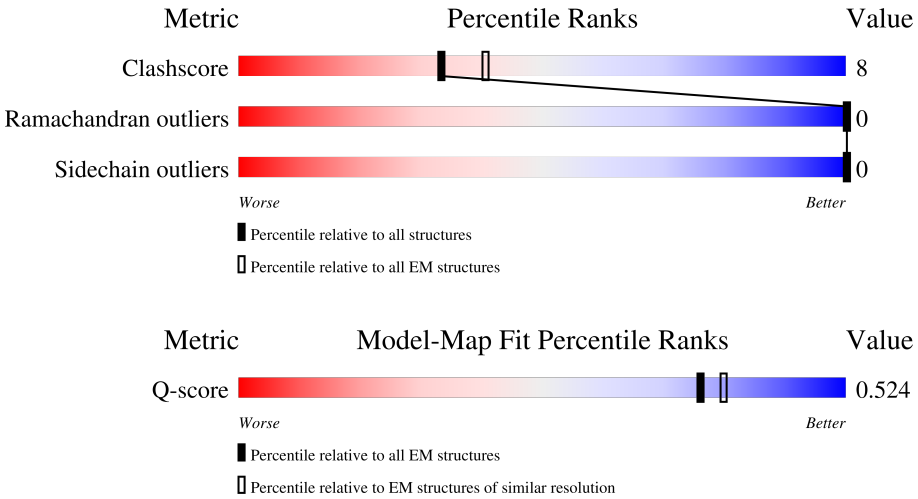
EMDB validation analysis	:	0.0.1.dev132
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.30 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	15087 (2.80 - 3.80)

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	396	12% · 85%
1	B	396	12% · 85%
1	C	396	12% · 85%
1	D	396	12% · 85%

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Mol	Chain	Length	Quality of chain
1	E	396	
1	F	396	
1	G	396	
1	H	396	
1	I	396	
1	J	396	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 4840 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 Apolipoprotein A-IV.

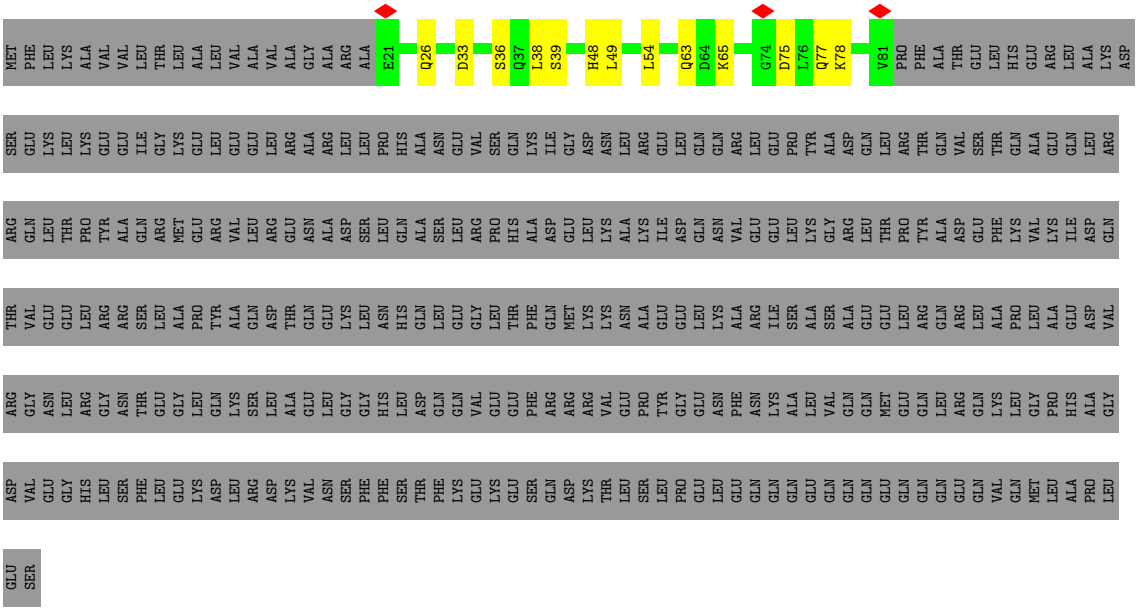
Mol	Chain	Residues	Atoms					AltConf	Trace
1	F	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	E	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	B	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	A	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	C	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	D	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	G	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	H	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	I	61	Total	C	N	O	S	0	0
			484	304	80	99	1		
1	J	61	Total	C	N	O	S	0	0
			484	304	80	99	1		

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.

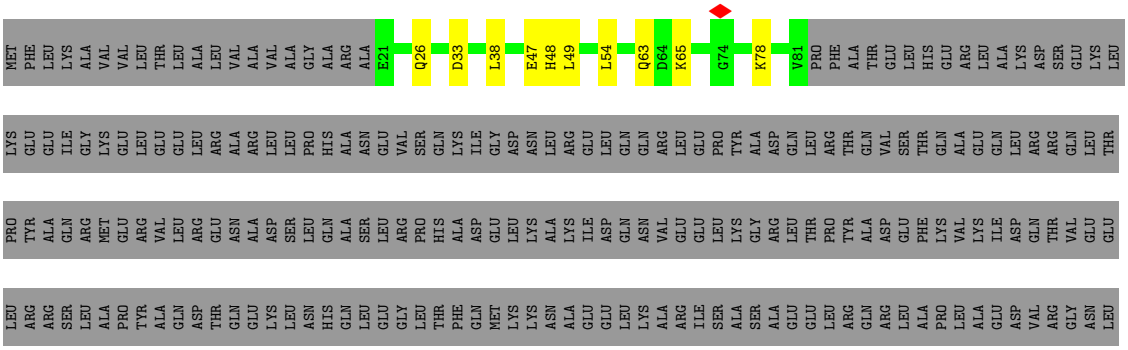
• Molecule 1: Apolipoprotein A-IV

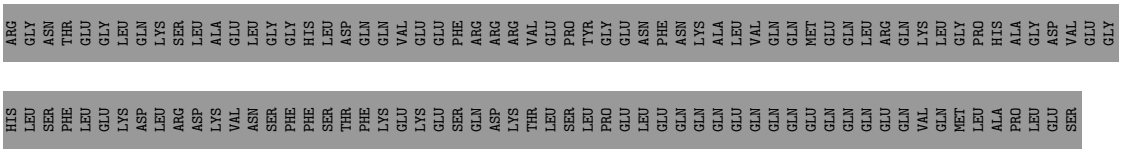
Chain F:



• Molecule 1: Apolipoprotein A-IV

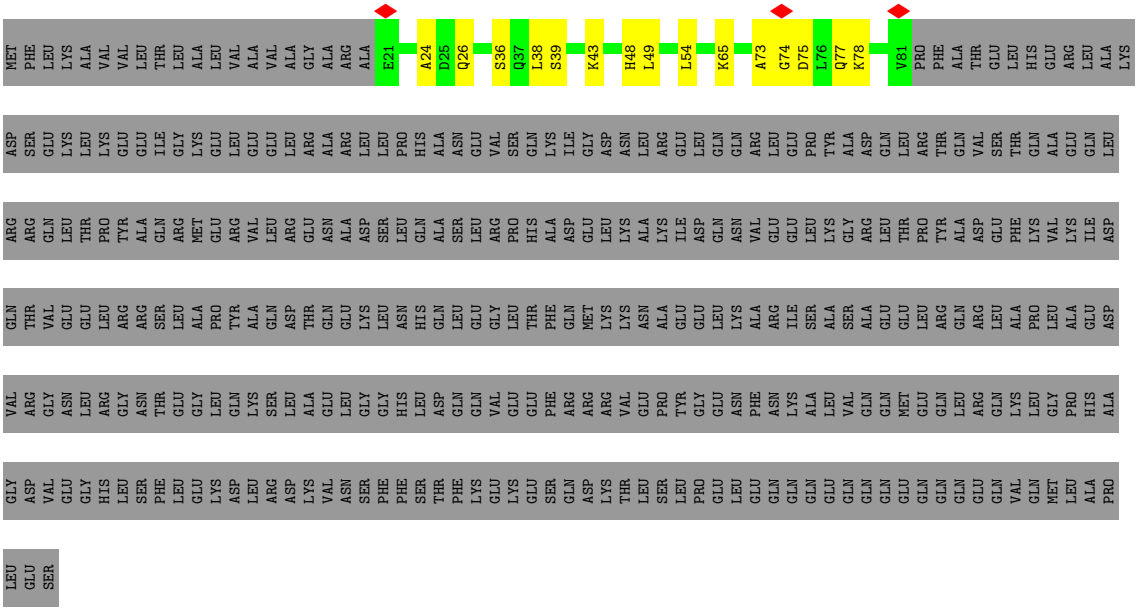
Chain E:





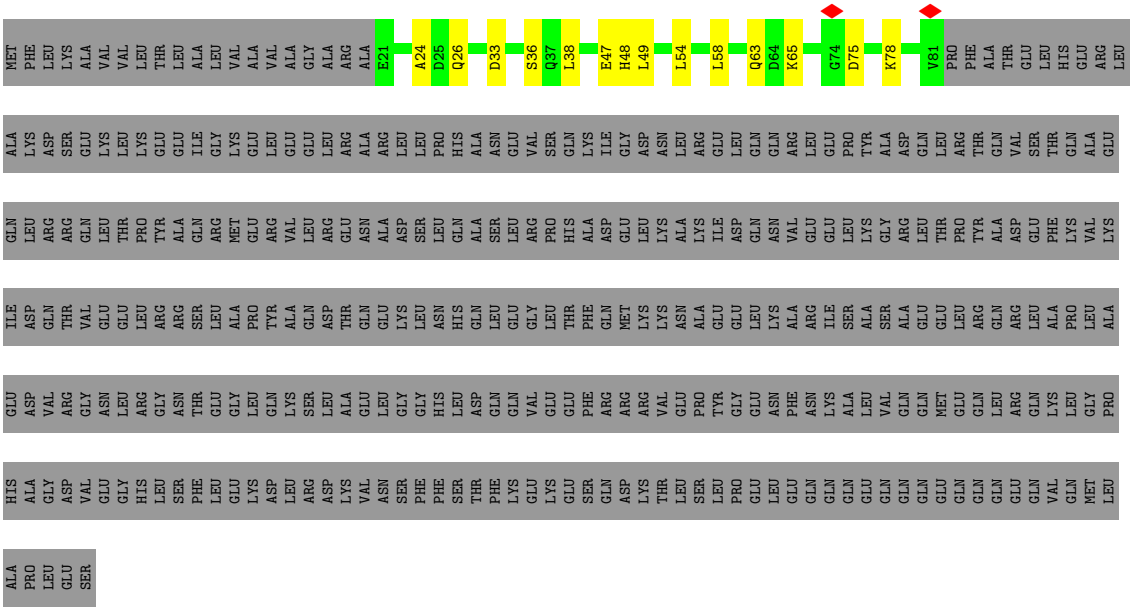
• Molecule 1: Apolipoprotein A-IV

Chain B:



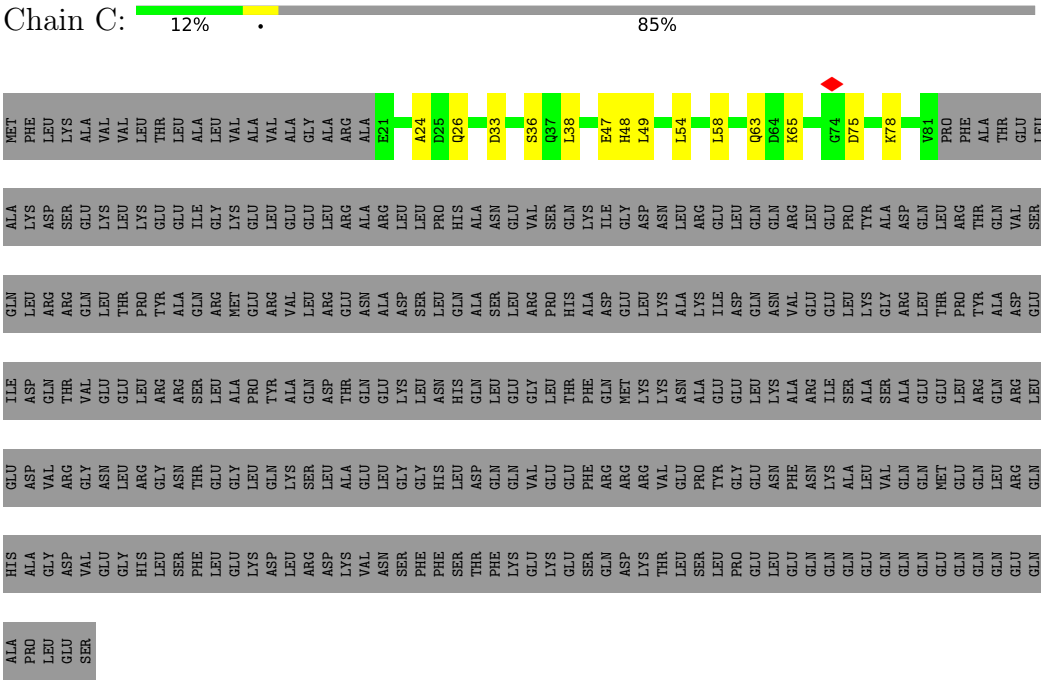
• Molecule 1: Apolipoprotein A-IV

Chain A:



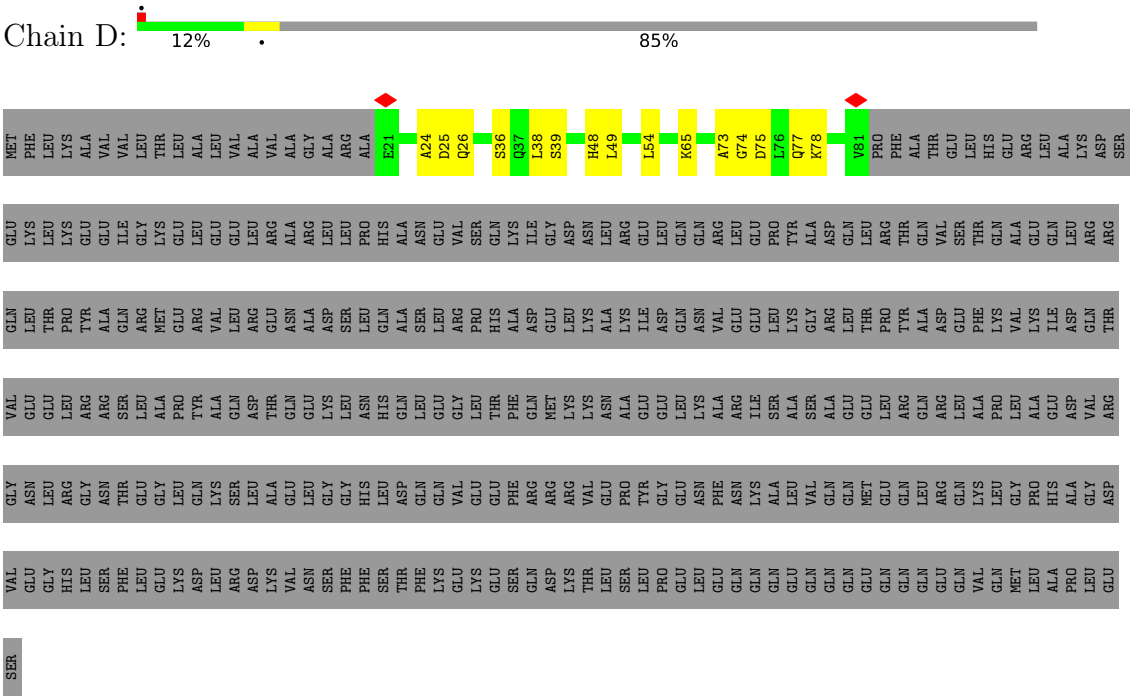
• Molecule 1: Apolipoprotein A-IV

Chain C:



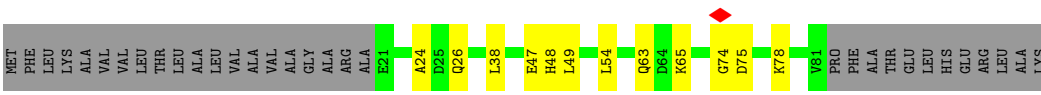
• Molecule 1: Apolipoprotein A-IV

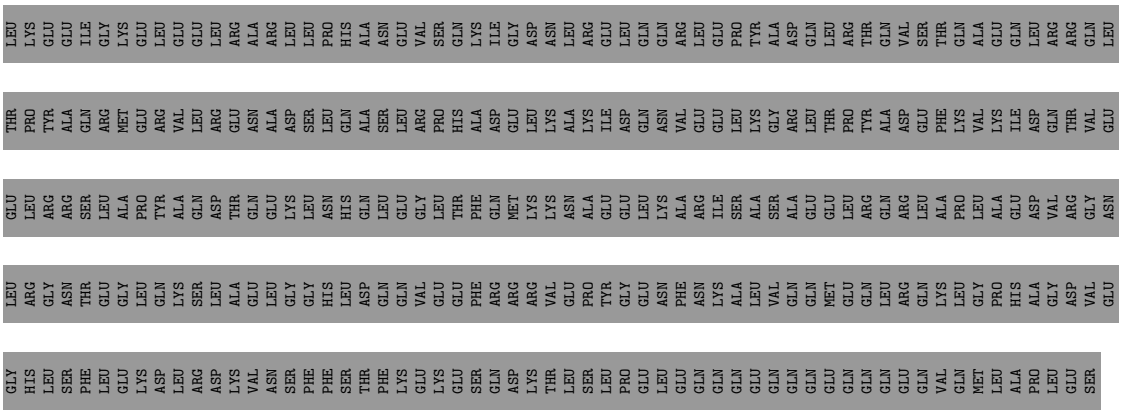
Chain D:



• Molecule 1: Apolipoprotein A-IV

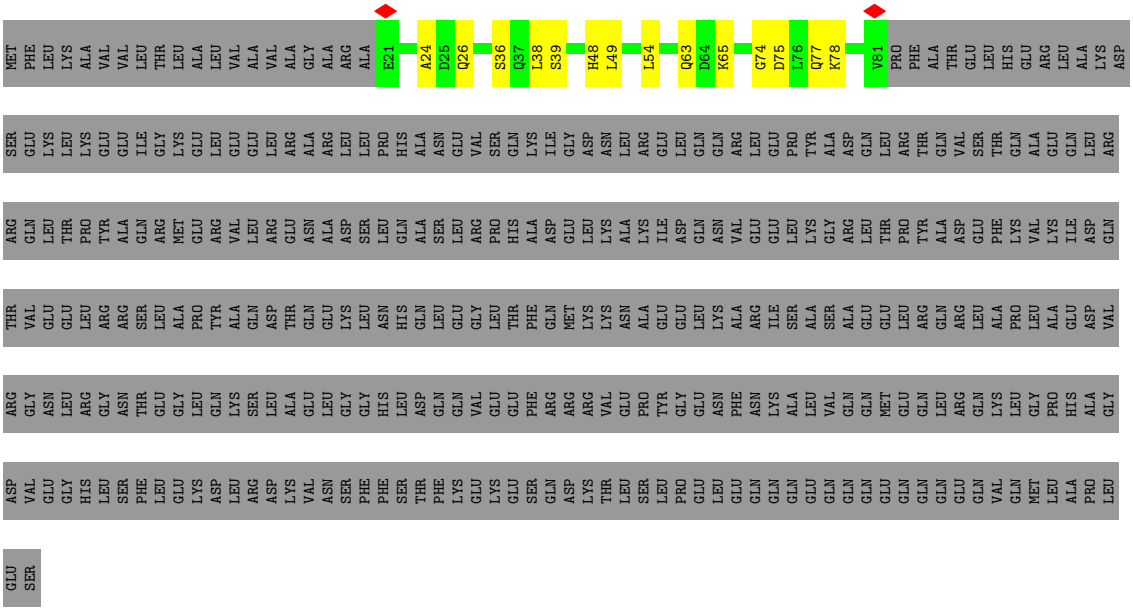
Chain G:





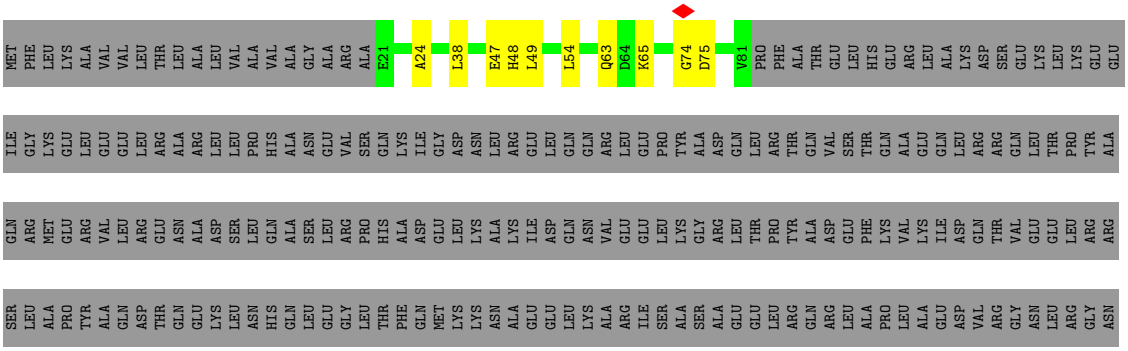
• Molecule 1: Apolipoprotein A-IV

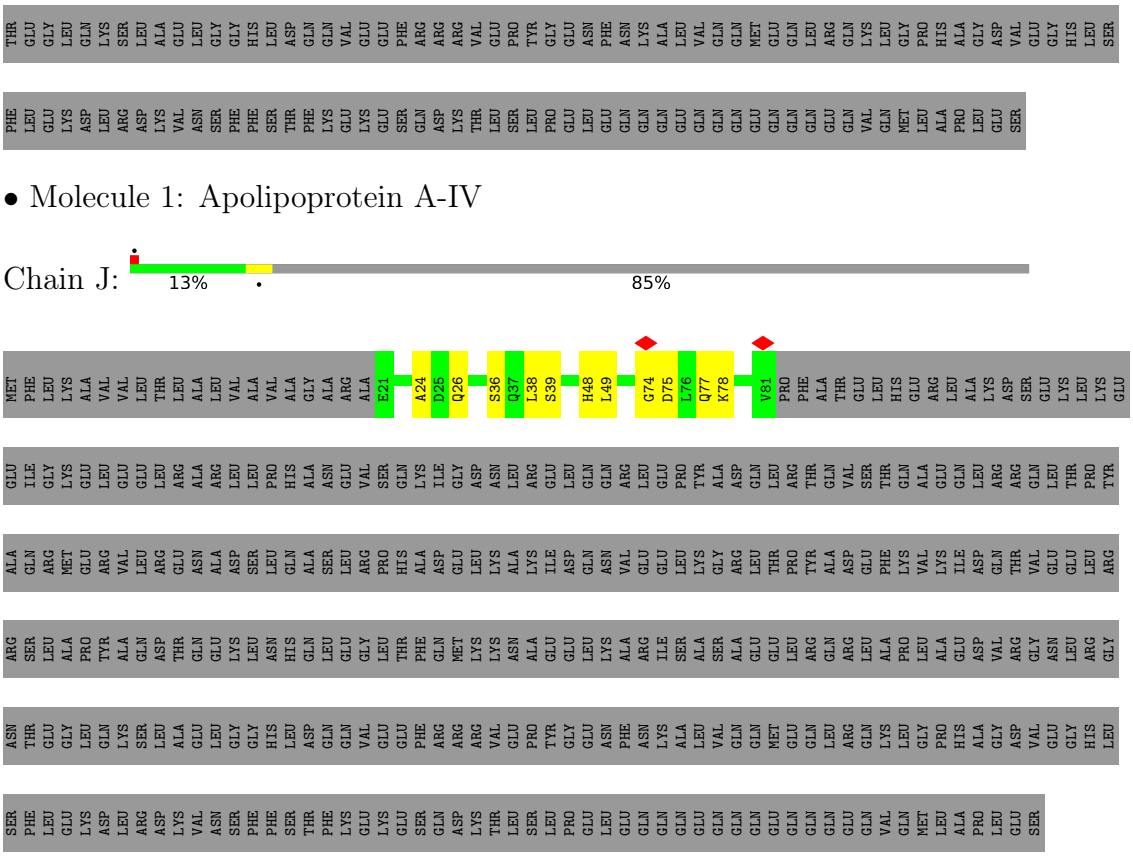
Chain H:



• Molecule 1: Apolipoprotein A-IV

Chain I:




THR GLU GLY LEU GLN LYS SER LEU ALA GLU LEU GLY GLY HIS LEU ASP GLN GLN VAL GLU GLU PHE ARG ARG ARG VAL GLU PRO TYR GLY GLU ASN PHE ASN LYS ALA LEU VAL GLN GLN MET GLU GLN LEU ARG GLN LYS LEU GLY PRO HIS ALA GLY ASP VAL GLU GLY HIS LEU SER
PHE LEU GLU LYS ASP LEU ARG ASP LYS VAL ASN SER PHE PHE SER THR PHE LYS GLU LYS GLU SER GLN ASP LYS THR LEU SER LEU PRO GLU LEU GLU GLN GLN GLN GLU GLN GLN GLN GLN GLU GLN GLN GLN GLU GLN VAL GLN MET LEU ALA PRO LEU GLU SER
Molecule 1: Apolipoprotein A-IV
Chain J:
13%
85%
MET PHE LEU LYS ALA VAL VAL LEU THR LEU ALA ALA LEU VAL ALA VAL ALA GLY ALA ARG ALA E21 A24 P25 Q26 S36 Q37 L38 S39 H48 L49 G74 D75 L76 Q77 K78 V81 PRO PHE ALA THR GLU LEU HIS GLU ARG LEU ALA LYS ASP SER GLU LYS LEU LYS GLU
GLU ILE GLY LYS GLU LEU GLU GLU LEU ARG ALA ARG LEU LEU PRO HIS ALA ASN GLU VAL SER GLN LYS ILE GLY ASP ASN LEU ARG GLU LEU GLN GLN ARG LEU GLU PRO TYR ALA ASP GLN LEU ARG THR GLN VAL SER THR GLN ALA GLU GLN LEU ARG ARG GLN LEU THR PRO TYR
ALA GLN ARG MET GLU ARG VAL LEU ARG GLU ASN ALA ASP SER LEU GLN ALA SER LEU ARG PRO HIS ALA ASP GLU LEU LYS ALA LYS ILE ASP GLN ASN VAL GLU GLU LEU LYS GLY ARG LEU THR PRO TYR ALA ASP GLU PHE LYS VAL LYS ILE ASP GLN THR VAL GLU GLU LEU ARG
ARG SER LEU ALA PRO TYR ALA GLN ASP THR GLN GLN GLU LYS LEU ASN HIS GLN LEU GLU GLY LEU THR PHE GLN MET LYS LYS ASN ALA GLU GLU LEU LYS ALA ARG ILE SER ALA SER ALA GLU GLU LEU ARG GLN ARG LEU ALA PRO LEU ALA GLU ASP VAL ARG GLY ASN LEU ARG GLY
ASN THR GLU GLY LEU GLN LYS SER LEU ALA GLU LEU GLY GLY HIS LEU ASP GLN GLN VAL GLU GLU PHE ARG ARG ARG VAL GLU PRO TYR GLY GLU ASN PHE ASN LYS ALA LEU VAL GLN GLN MET GLU GLN LEU ARG GLN LYS LEU GLY PRO HIS ALA GLY ASP VAL GLU GLY HIS LEU
SER PHE LEU GLU LYS ASP LEU ARG ASP LYS VAL ASN SER PHE PHE SER THR PHE LYS GLU LYS GLU SER GLN ASP LYS THR LEU SER LEU PRO GLU LEU GLU GLN GLN GLN GLU GLN GLN GLN GLN GLU GLN VAL GLN MET LEU ALA PRO LEU GLU SER

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-1.04°, rise=4.76 Å, axial sym=C1	Depositor
Number of segments used	36124	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 ($e^-/Å^2$)	60	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.053	Depositor
Minimum map value	-0.028	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	230.4, 230.4, 230.4	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.28, 1.28, 1.28	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	0.10	0/490	0.30	0/662
1	B	0.11	0/490	0.33	0/662
1	C	0.10	0/490	0.30	0/662
1	D	0.11	0/490	0.33	0/662
1	E	0.10	0/490	0.30	0/662
1	F	0.11	0/490	0.33	0/662
1	G	0.10	0/490	0.30	0/662
1	H	0.11	0/490	0.33	0/662
1	I	0.10	0/490	0.30	0/662
1	J	0.11	0/490	0.33	0/662
All	All	0.10	0/4900	0.31	0/6620

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	484	0	471	12	0
1	B	484	0	471	13	0
1	C	484	0	471	16	0
1	D	484	0	471	15	0
1	E	484	0	471	11	0
1	F	484	0	471	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	484	0	471	14	0
1	H	484	0	471	12	0
1	I	484	0	471	11	0
1	J	484	0	471	8	0
All	All	4840	0	4710	79	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 8.

All (79) 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:E:33:ASP:OD2	1:I:65:LYS:NZ	2.23	0.71
1:E:65:LYS:NZ	1:A:33:ASP:OD2	2.25	0.68
1:H:38:LEU:HD13	1:J:38:LEU:HD23	1.76	0.67
1:A:38:LEU:HD23	1:C:38:LEU:HD13	1.77	0.66
1:A:54:LEU:HD12	1:C:54:LEU:HD22	1.78	0.66
1:C:33:ASP:OD2	1:G:65:LYS:NZ	2.28	0.65
1:F:38:LEU:HD23	1:D:38:LEU:HD13	1.80	0.64
1:F:38:LEU:HD13	1:H:38:LEU:HD23	1.81	0.63
1:G:38:LEU:HD23	1:I:38:LEU:HD13	1.80	0.63
1:G:54:LEU:HD12	1:I:54:LEU:HD22	1.80	0.63
1:B:38:LEU:HD13	1:D:38:LEU:HD23	1.83	0.61
1:E:38:LEU:HD13	1:C:38:LEU:HD23	1.82	0.60
1:D:36:SER:HG	1:D:39:SER:HG	1.47	0.60
1:E:54:LEU:HD12	1:G:54:LEU:HD22	1.83	0.60
1:E:38:LEU:HD23	1:G:38:LEU:HD13	1.84	0.59
1:E:54:LEU:HD22	1:C:54:LEU:HD12	1.85	0.59
1:F:36:SER:OG	1:F:39:SER:OG	2.21	0.54
1:H:36:SER:OG	1:H:39:SER:OG	2.21	0.53
1:J:36:SER:OG	1:J:39:SER:OG	2.21	0.52
1:A:63:GLN:HB2	1:C:63:GLN:HG2	1.95	0.48
1:D:36:SER:OG	1:D:39:SER:OG	2.21	0.47
1:C:48:HIS:C	1:C:49:LEU:HD12	2.39	0.47
1:E:48:HIS:C	1:E:49:LEU:HD12	2.39	0.47
1:A:48:HIS:C	1:A:49:LEU:HD12	2.39	0.47
1:G:48:HIS:C	1:G:49:LEU:HD12	2.39	0.47
1:I:48:HIS:C	1:I:49:LEU:HD12	2.39	0.47
1:H:75:ASP:OD1	1:J:75:ASP:HB3	2.15	0.47
1:B:48:HIS:C	1:B:49:LEU:HD12	2.41	0.46
1:D:48:HIS:C	1:D:49:LEU:HD12	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ASP:HB3	1:C:75:ASP:OD1	2.16	0.46
1:D:26:GLN:NE2	1:D:78:LYS:O	2.49	0.45
1:B:26:GLN:NE2	1:B:78:LYS:O	2.49	0.45
1:B:75:ASP:OD1	1:D:75:ASP:HB3	2.16	0.45
1:F:48:HIS:C	1:F:49:LEU:HD12	2.41	0.45
1:B:24:ALA:HA	1:D:24:ALA:O	2.16	0.45
1:G:75:ASP:HB3	1:I:75:ASP:OD1	2.17	0.45
1:A:65:LYS:O	1:C:65:LYS:HA	2.16	0.45
1:G:65:LYS:O	1:I:65:LYS:HA	2.16	0.45
1:H:48:HIS:C	1:H:49:LEU:HD12	2.41	0.45
1:A:58:LEU:HA	1:C:58:LEU:O	2.17	0.45
1:J:48:HIS:C	1:J:49:LEU:HD12	2.41	0.45
1:G:74:GLY:O	1:I:74:GLY:HA3	2.17	0.44
1:B:54:LEU:HD22	1:D:54:LEU:HD12	1.99	0.44
1:A:24:ALA:O	1:C:24:ALA:HA	2.18	0.44
1:G:24:ALA:O	1:I:24:ALA:HA	2.18	0.44
1:F:75:ASP:HB3	1:D:75:ASP:OD1	2.19	0.43
1:B:73:ALA:HA	1:D:73:ALA:O	2.19	0.43
1:J:26:GLN:NE2	1:J:78:LYS:O	2.49	0.43
1:B:65:LYS:HA	1:D:65:LYS:O	2.17	0.43
1:F:54:LEU:HD22	1:H:54:LEU:HD12	2.00	0.43
1:G:63:GLN:HB2	1:I:63:GLN:HG2	2.02	0.42
1:B:36:SER:OG	1:B:39:SER:OG	2.21	0.42
1:H:26:GLN:NE2	1:H:78:LYS:O	2.49	0.42
1:F:65:LYS:HA	1:H:65:LYS:O	2.19	0.42
1:E:63:GLN:HG2	1:C:63:GLN:HB2	2.02	0.42
1:B:74:GLY:HA3	1:D:74:GLY:O	2.19	0.42
1:I:38:LEU:HD11	1:I:47:GLU:CD	2.45	0.42
1:B:43:LYS:NZ	1:D:25:ASP:OD2	2.47	0.42
1:G:38:LEU:HD11	1:G:47:GLU:CD	2.45	0.41
1:F:26:GLN:NE2	1:F:78:LYS:O	2.49	0.41
1:E:38:LEU:HD11	1:E:47:GLU:CD	2.45	0.41
1:B:26:GLN:O	1:B:77:GLN:NE2	2.53	0.41
1:A:38:LEU:HD11	1:A:47:GLU:CD	2.45	0.41
1:C:38:LEU:HD11	1:C:47:GLU:CD	2.45	0.41
1:D:26:GLN:O	1:D:77:GLN:NE2	2.53	0.41
1:F:33:ASP:OD2	1:B:65:LYS:NZ	2.37	0.41
1:A:26:GLN:NE2	1:A:78:LYS:O	2.51	0.41
1:F:26:GLN:O	1:F:77:GLN:NE2	2.53	0.41
1:J:26:GLN:O	1:J:77:GLN:NE2	2.53	0.41
1:E:65:LYS:HA	1:C:65:LYS:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:GLN:NE2	1:C:78:LYS:O	2.51	0.41
1:E:26:GLN:NE2	1:E:78:LYS:O	2.51	0.41
1:H:74:GLY:HA3	1:J:74:GLY:O	2.21	0.41
1:F:63:GLN:HG2	1:H:63:GLN:HB2	2.02	0.40
1:A:36:SER:CB	1:C:36:SER:O	2.69	0.40
1:G:26:GLN:NE2	1:G:78:LYS:O	2.51	0.40
1:H:26:GLN:O	1:H:77:GLN:NE2	2.53	0.40
1:G:75:ASP:CB	1:I:75:ASP:OD1	2.70	0.40
1:H:24:ALA:HA	1:J:24:ALA:O	2.20	0.40

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	59/396 (15%)	56 (95%)	3 (5%)	0	100	100
1	B	59/396 (15%)	55 (93%)	4 (7%)	0	100	100
1	C	59/396 (15%)	56 (95%)	3 (5%)	0	100	100
1	D	59/396 (15%)	55 (93%)	4 (7%)	0	100	100
1	E	59/396 (15%)	56 (95%)	3 (5%)	0	100	100
1	F	59/396 (15%)	55 (93%)	4 (7%)	0	100	100
1	G	59/396 (15%)	56 (95%)	3 (5%)	0	100	100
1	H	59/396 (15%)	55 (93%)	4 (7%)	0	100	100
1	I	59/396 (15%)	56 (95%)	3 (5%)	0	100	100
1	J	59/396 (15%)	55 (93%)	4 (7%)	0	100	100
All	All	590/3960 (15%)	555 (94%)	35 (6%)	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	53/344 (15%)	53 (100%)	0	100	100
1	B	53/344 (15%)	53 (100%)	0	100	100
1	C	53/344 (15%)	53 (100%)	0	100	100
1	D	53/344 (15%)	53 (100%)	0	100	100
1	E	53/344 (15%)	53 (100%)	0	100	100
1	F	53/344 (15%)	53 (100%)	0	100	100
1	G	53/344 (15%)	53 (100%)	0	100	100
1	H	53/344 (15%)	53 (100%)	0	100	100
1	I	53/344 (15%)	53 (100%)	0	100	100
1	J	53/344 (15%)	53 (100%)	0	100	100
All	All	530/3440 (15%)	530 (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 (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	63	GLN
1	A	63	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.

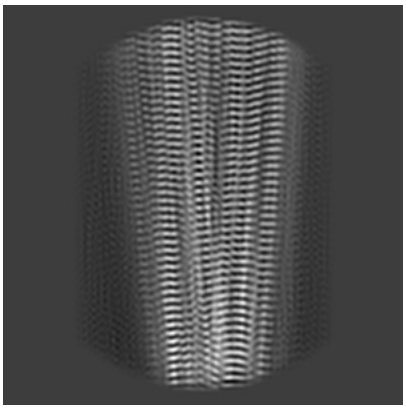
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-52457. These allow visual inspection of the internal detail of the map and identification of artifacts.

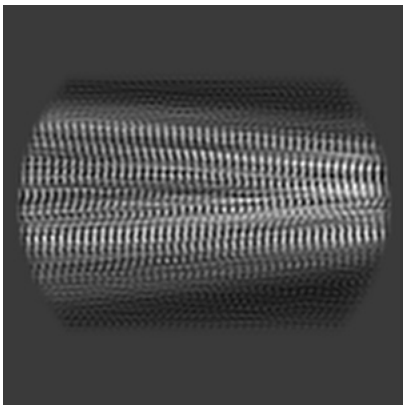
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

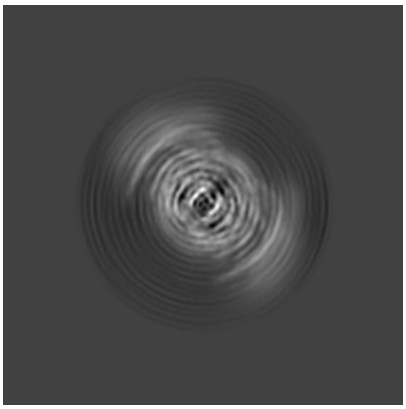
6.1.1 Primary map



X

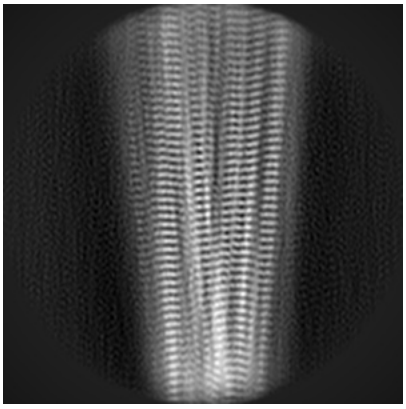


Y

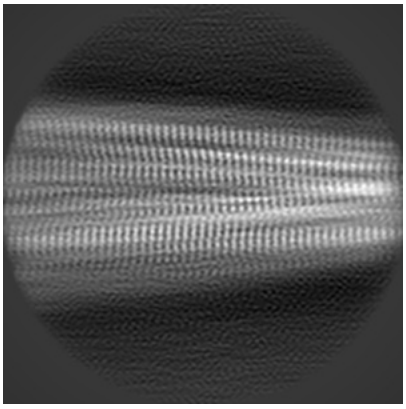


Z

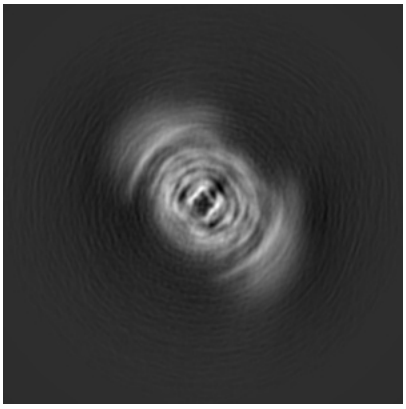
6.1.2 Raw map



X



Y

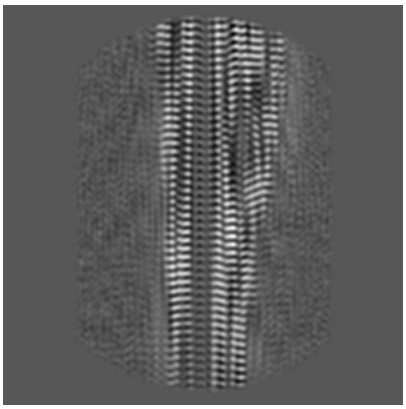


Z

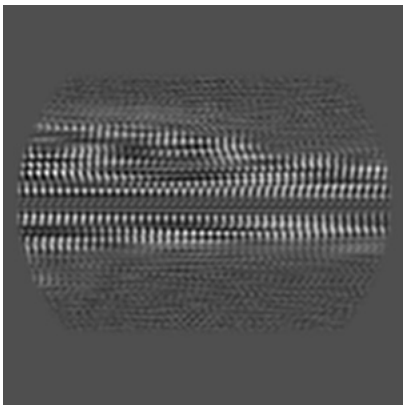
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

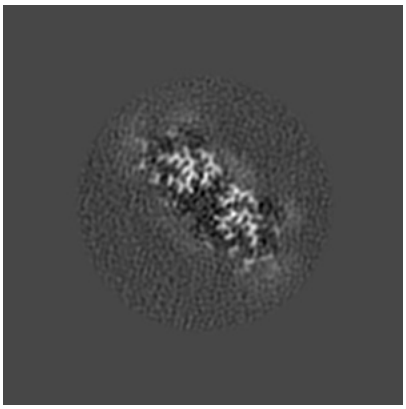
6.2.1 Primary map



X Index: 90

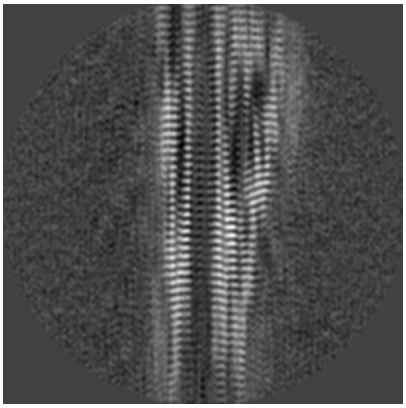


Y Index: 90

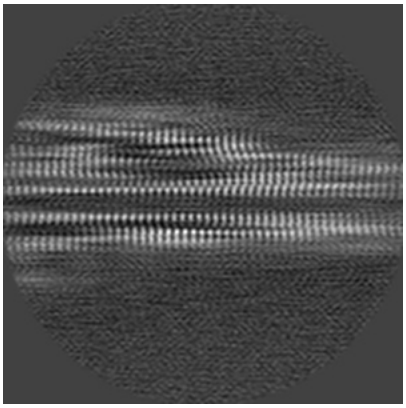


Z Index: 90

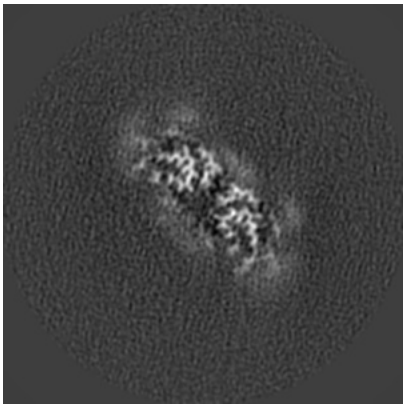
6.2.2 Raw map



X Index: 90



Y Index: 90

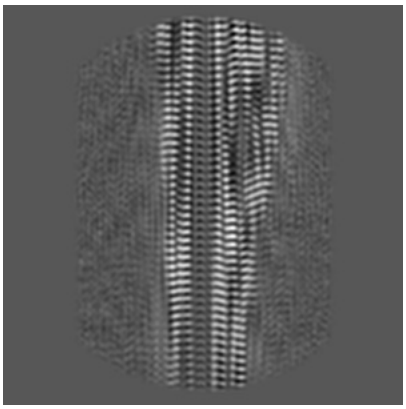


Z Index: 90

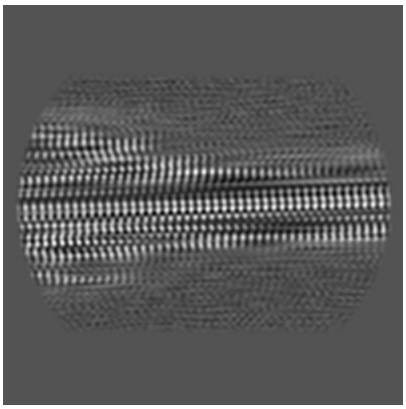
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

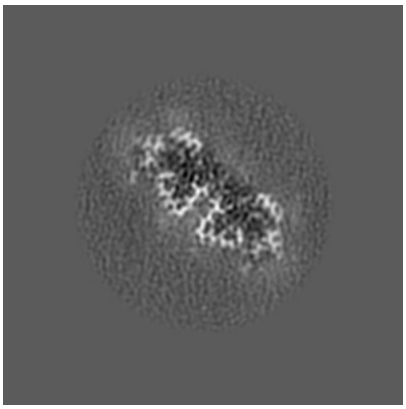
6.3.1 Primary map



X Index: 90

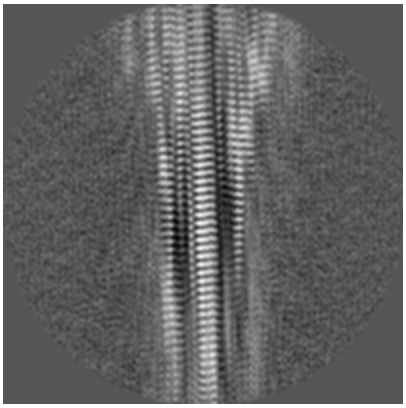


Y Index: 96

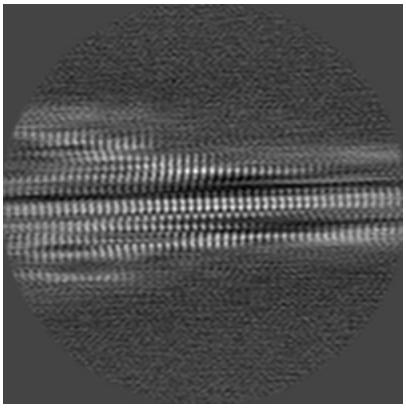


Z Index: 77

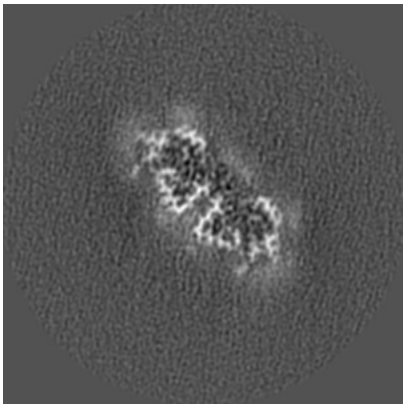
6.3.2 Raw map



X Index: 96



Y Index: 96

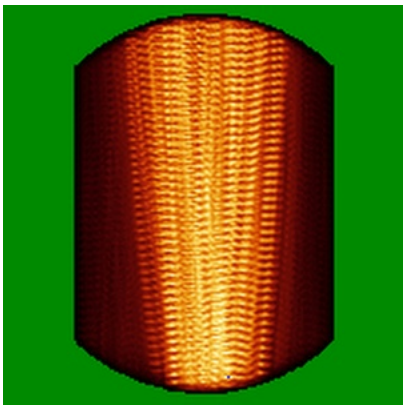


Z Index: 92

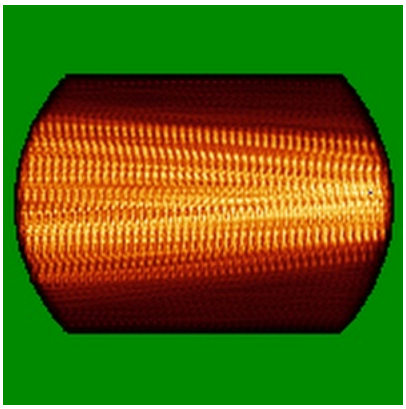
The images above show the largest variance slices of the map in three orthogonal directions.

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

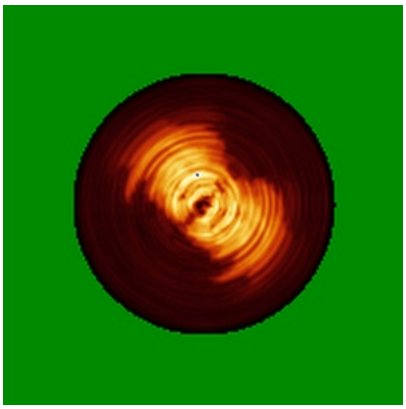
6.4.1 Primary map



X

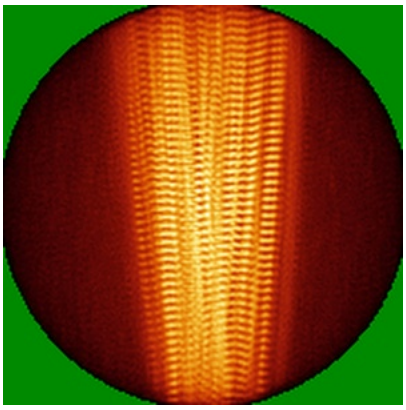


Y

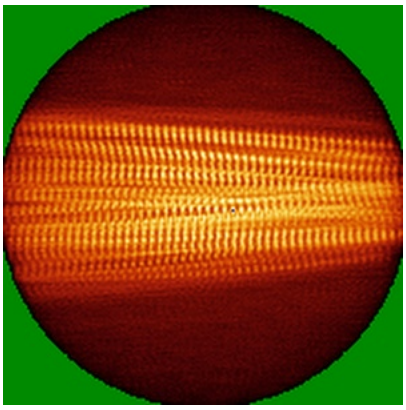


Z

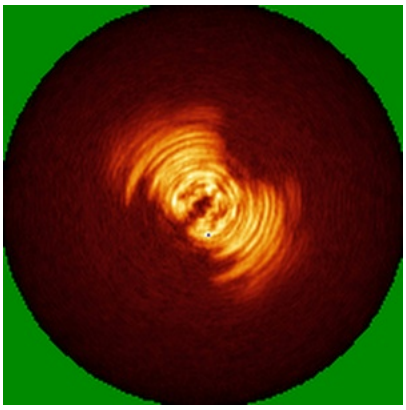
6.4.2 Raw map



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)

This section was not generated.

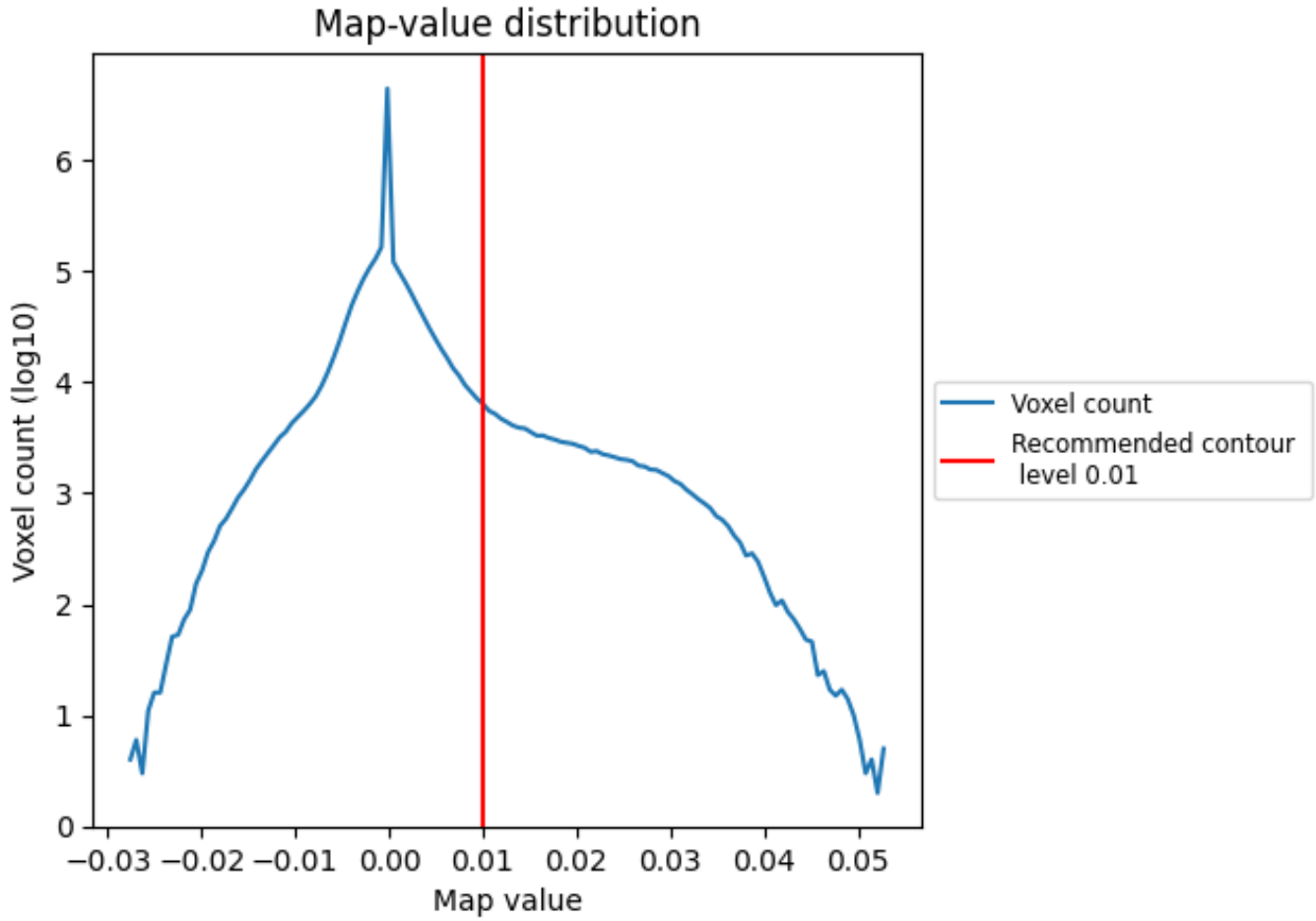
6.6 Mask visualisation (i)

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

7 Map analysis

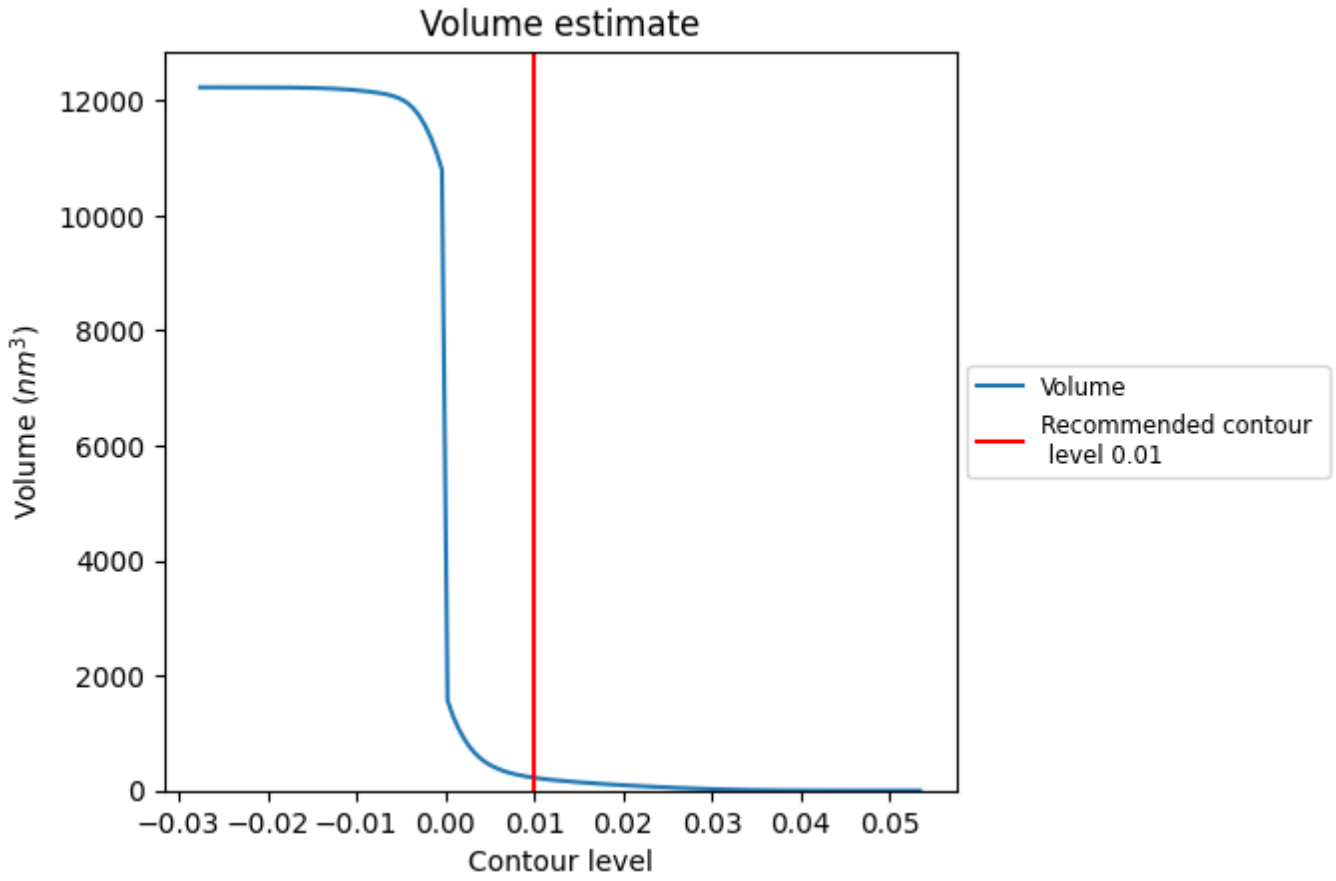
This section contains the results of statistical analysis of the map.

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.

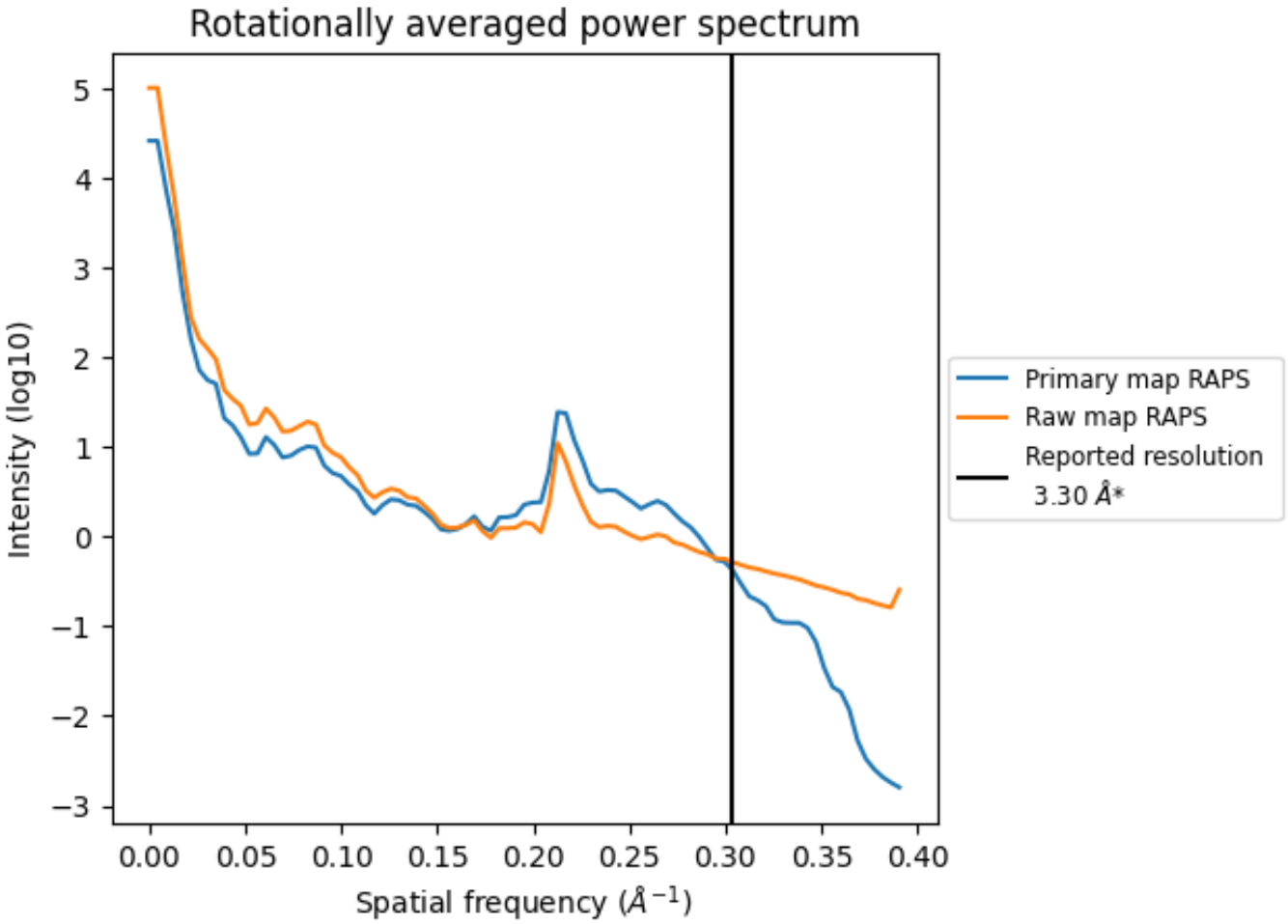
7.2 Volume estimate



The volume at the recommended contour level is 223 nm^3; this corresponds to an approximate mass of 202 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)

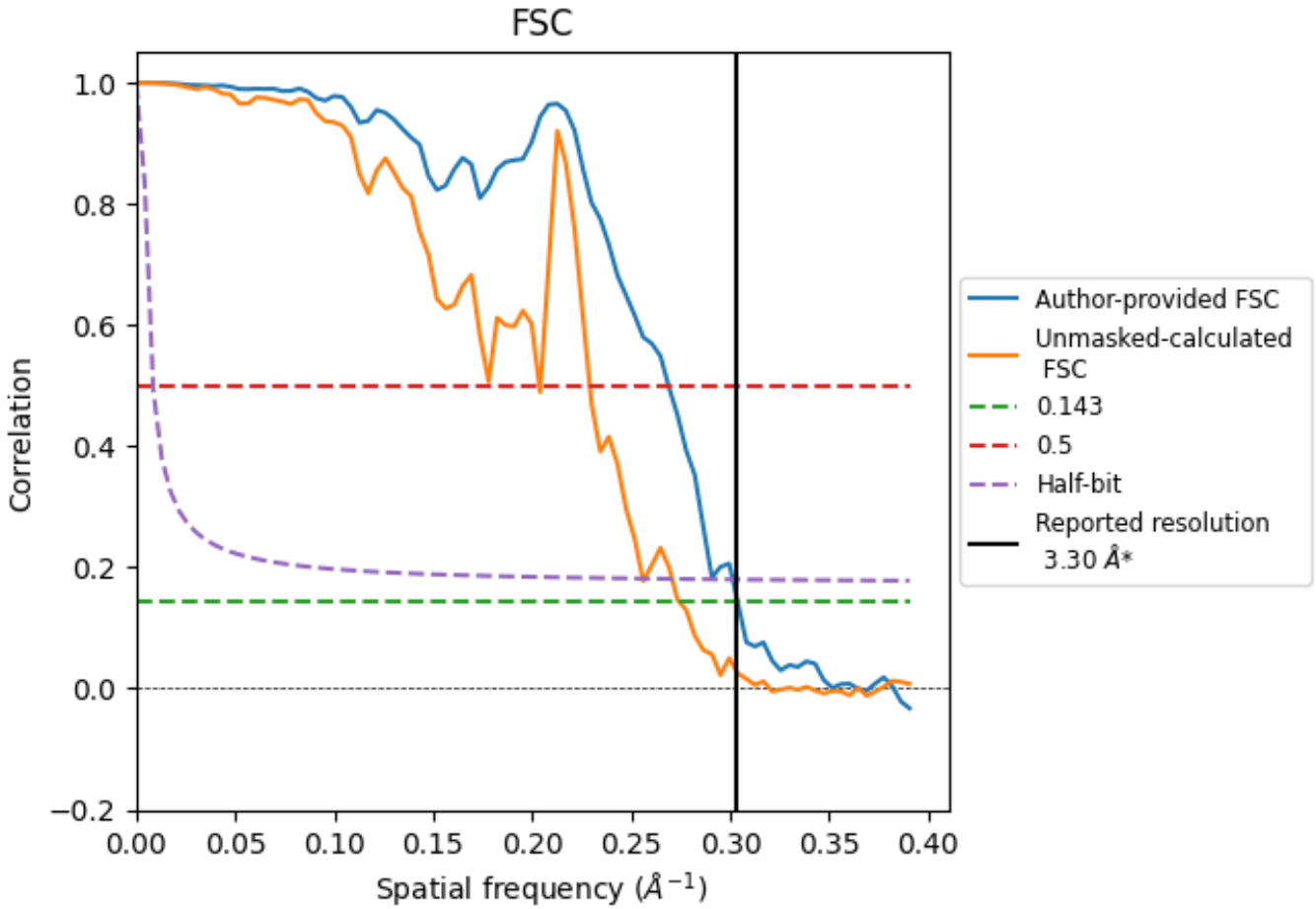


*Reported resolution corresponds to spatial frequency of 0.303 Å^{-1}

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 0.303 Å^{-1}

8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.29	3.72	3.32
Unmasked-calculated*	3.65	4.91	3.91

*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 3.65 differs from the reported value 3.3 by more than 10 %

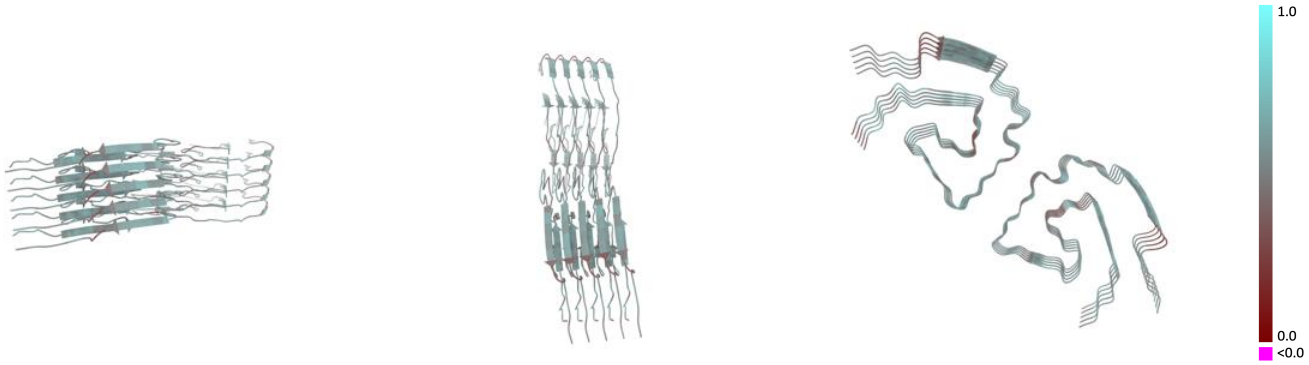
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-52457 and PDB model 9HX4. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay

This section was not generated.

9.2 Q-score mapped to coordinate model



X Y Z

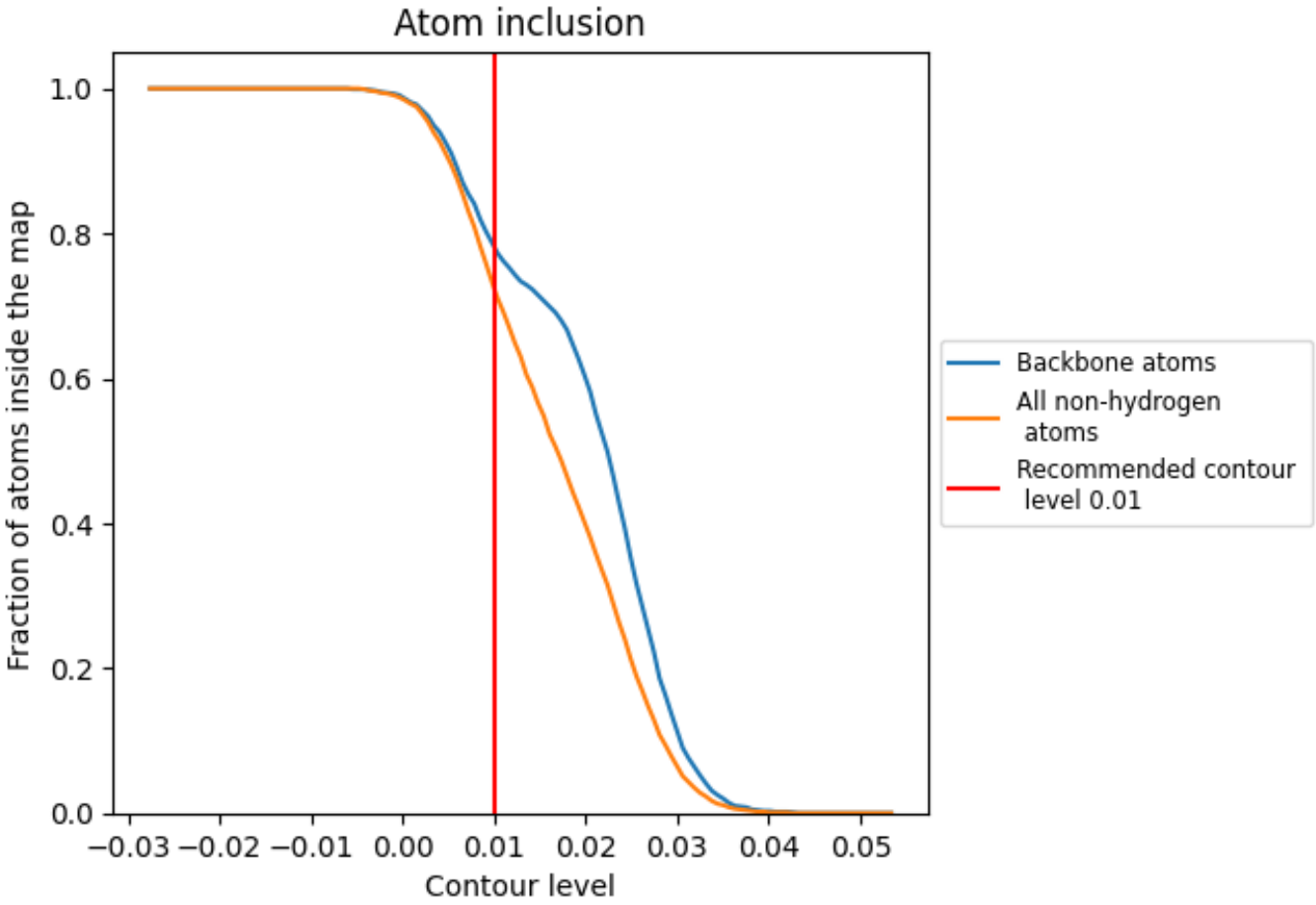
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).

9.4 Atom inclusion (i)



At the recommended contour level, 78% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7240	0.5240
A	0.7170	0.5200
B	0.7340	0.5270
C	0.7130	0.5230
D	0.7240	0.5260
E	0.7240	0.5240
F	0.7340	0.5230
G	0.7150	0.5250
H	0.7340	0.5260
I	0.7190	0.5250
J	0.7260	0.5200