



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 01:34 PM UTC

PDB ID : 7NCH / pdb_00007nch
EMDB ID : EMD-12266
Title : Type 1B alpha-synuclein filament seeded in vitro by filaments purified from Multiple Systems Atrophy Case 1
Authors : Lovestam, S.K.A.; Schweighauser, M.; Scheres, S.H.W.
Deposited on : 2021-01-28
Resolution : 3.84 Å(reported)

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

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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:

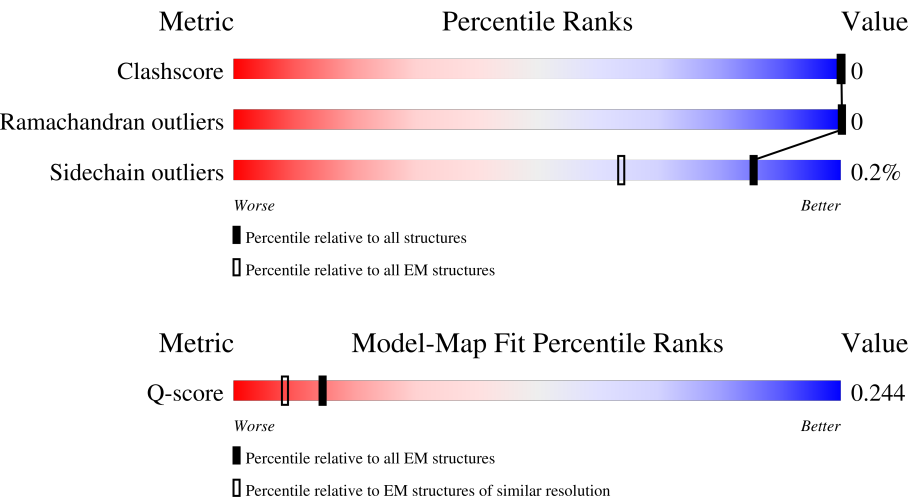
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.84 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.



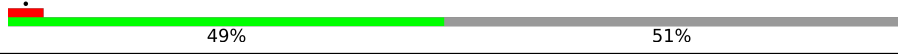
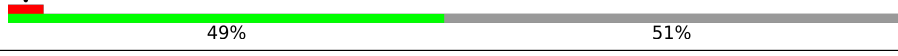
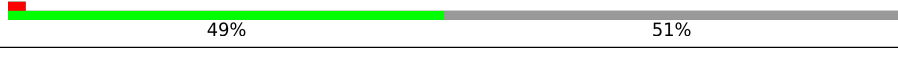
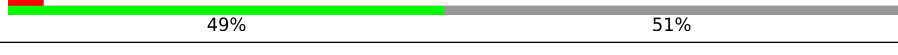
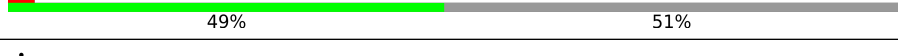
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	9033 (3.34 - 4.34)

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	140	49% 51%
1	B	140	49% 51%
1	C	140	49% 51%
1	D	140	48% 51%

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Mol	Chain	Length	Quality of chain
1	E	140	 49%51%
1	F	140	 49%51%
1	G	140	 49%51%
1	H	140	 49%51%
1	I	140	 49%51%
1	J	140	 49%51%
1	K	140	 49%51%
1	L	140	 49%51%

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 11292 atoms, of which 5796 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 Alpha-synuclein.

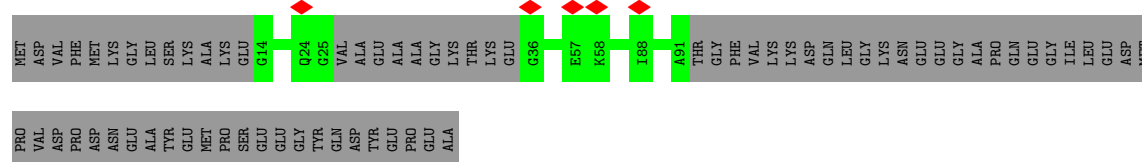
Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	B	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	C	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	D	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	E	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	F	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	G	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	H	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	I	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	J	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	K	68	Total 941	C 284	H 483	N 81	O 93	0	0
1	L	68	Total 941	C 284	H 483	N 81	O 93	0	0

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.

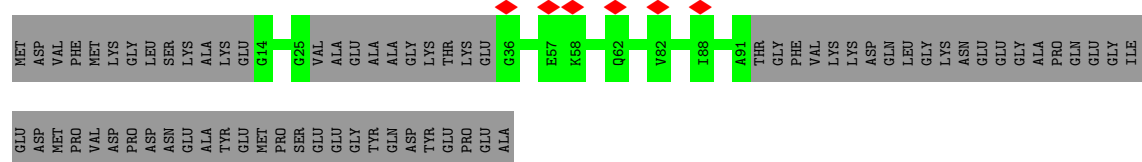
- Molecule 1: Alpha-synuclein

Chain A: 



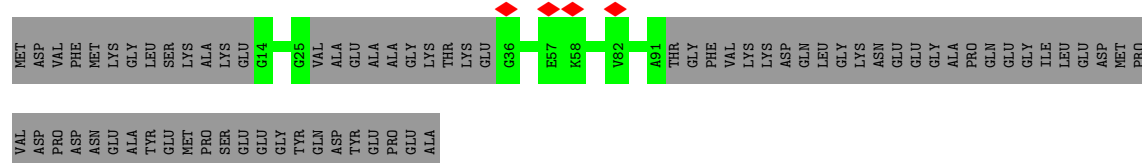
- Molecule 1: Alpha-synuclein

Chain B: 



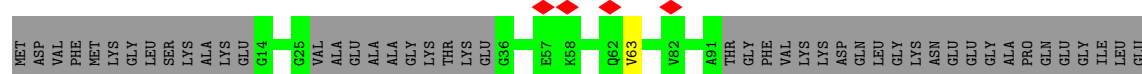
- Molecule 1: Alpha-synuclein

Chain C: 



- Molecule 1: Alpha-synuclein

Chain D: 



ASP
MET
PRO
VAL
PHE
ASP
PRO
LYS
ASP
ASN
GLU
TYR
TRP
GLY
MET
PRO
SER
GLU
GLY
TYR
GLN
ASP
TYR
GLU
PRO
GLU
ALA

• Molecule 1: Alpha-synuclein

Chain E:  49% 51%

MET
ASP
VAL
PHE
ASP
MET
LYS
GLY
LEU
SER
LYS
ALA
LYS
LYS
GLU
G14
G25
VAL
ALA
GLY
GLY
ALA
ALA
GLY
LYS
THR
LYS
GLU
G36
E57
K58
V62
A91
THR
GLY
PHE
VAL
LYS
LYS
ASP
GLN
LEU
GLY
LYS
ASN
GLY
GLU
GLY
ALA
PRO
GLN
GLY
ILE
LEU
GLU
ASP
MET
PRO

VAL
ASP
PRO
ASP
ASN
GLY
ALA
TYR
GLU
MET
PRO
LYS
SER
GLU
GLY
TYR
GLN
ASP
TYR
GLU
PRO
ALA

• Molecule 1: Alpha-synuclein

Chain F:  49% 51%

MET
ASP
VAL
PHE
ASP
MET
LYS
GLY
LEU
SER
LYS
ALA
LYS
LYS
GLU
G14
G25
VAL
ALA
GLY
GLY
ALA
ALA
GLY
LYS
THR
LYS
GLU
G36
E57
K58
Q62
V62
A91
THR
GLY
PHE
VAL
LYS
LYS
ASP
GLN
LEU
LYS
ASN
GLY
GLU
GLY
ALA
PRO
GLN
GLY
ILE
LEU
GLU
ASP

MET
PRO
VAL
ASP
PRO
ASP
ASN
GLU
ALA
TYR
GLU
MET
PRO
SER
GLU
GLY
TYR
GLN
ASP
TYR
GLU
PRO
GLU
ALA

• Molecule 1: Alpha-synuclein

Chain G:  49% 51%

MET
ASP
VAL
PHE
ASP
MET
LYS
GLY
LEU
SER
LYS
ALA
LYS
LYS
GLU
G14
G24
G25
VAL
ALA
GLY
GLY
ALA
ALA
GLY
LYS
THR
LYS
GLU
G36
E57
K58
V62
I88
A91
THR
GLY
PHE
VAL
LYS
LYS
ASP
GLN
LEU
LYS
ASN
GLY
GLU
GLY
ALA
PRO
GLN
GLY
ILE
LEU
GLU

ASP
MET
PRO
VAL
PHE
ASP
PRO
LYS
ASP
ASN
GLU
TYR
TRP
GLY
MET
PRO
SER
GLU
GLY
TYR
GLN
ASP
TYR
GLU
PRO
GLU
ALA

• Molecule 1: Alpha-synuclein

Chain H:  49% 51%

MET
ASP
VAL
PHE
ASP
MET
LYS
GLY
LEU
SER
LYS
ALA
LYS
LYS
GLU
G14
G25
VAL
ALA
GLY
GLY
ALA
ALA
GLY
LYS
THR
LYS
GLU
G36
E57
K58
Q62
V62
I88
A91
THR
GLY
PHE
VAL
LYS
LYS
ASP
GLN
LEU
GLY
LYS
ASN
GLY
GLU
GLY
ALA
PRO
GLN
GLY
ILE
LEU

GLU
ASP
MET
PRO
VAL
PHE
ASP
PRO
LYS
ASP
ASN
GLU
TYR
TRP
GLY
MET
PRO
SER
GLU
GLY
TYR
GLN
ASP
TYR
GLU
PRO
GLU
ALA

• Molecule 1: Alpha-synuclein

Chain I:  49% 51%

MET
ASP
VAL
PHE
ASP
MET
LYS
GLY
LEU
SER
LYS
ALA
LYS
LYS
GLU
G14
G25
VAL
ALA
GLY
GLY
ALA
ALA
GLY
LYS
THR
LYS
GLU
G36
E57
K58
V62
A91
THR
GLY
PHE
VAL
LYS
LYS
ASP
GLN
LEU
GLY
LYS
ASN
GLY
GLU
GLY
ALA
PRO
GLN
GLY
ILE
LEU
GLU
ASP
MET
PRO

VAL ASP
PRO PRO
ASP ASP
ASN ASN
GLU GLU
ALA ALA
TYR TYR
MET MET
PRO PRO
SER SER
GLU GLU
GLY GLY
GLN GLN
ASP ASP
TYR TYR
GLU GLU
PRO PRO
GLU GLU
ALA ALA

● Molecule 1: Alpha-synuclein



MET ASP
PRO VAL
ASP PHE
PRO MET
ASP LYS
ASN GLY
GLU LEU
TYR SER
LYS TYR
ALA ALA
MET LYS
GLU PRO
SER GLU
G14 G25
VAL VAL
ALA ALA
GLU GLU
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
GLU GLU
G36 E57 K58 Q62 V62 A91
THR THR
GLY GLY
PHE PHE
VAL VAL
LYS LYS
LYS LYS
ASP ASP
GLN GLN
LEU LEU
GLY GLY
LYS LYS
ASN ASN
GLU GLU
GLY GLY
ALA ALA
PRO PRO
GLN GLN
GLY GLY
ILE ILE
LEU LEU
GLU GLU
ASP ASP

MET PRO
VAL VAL
ASP ASP
PRO PRO
ASP ASP
ASN ASN
GLU GLU
TYR TYR
LYS LYS
ALA ALA
MET MET
PRO PRO
SER SER
GLU GLU
GLY GLY
TYR TYR
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
ALA ALA

● Molecule 1: Alpha-synuclein



MET ASP
PRO VAL
ASP PHE
PRO MET
ASP LYS
ASN GLY
GLU LEU
TYR SER
LYS TYR
ALA ALA
MET LYS
GLU PRO
SER GLU
G14 G25
VAL VAL
ALA ALA
GLU GLU
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
GLU GLU
G36 E57 K58 V62 I88 A91
THR THR
GLY GLY
PHE PHE
VAL VAL
LYS LYS
LYS LYS
ASP ASP
GLN GLN
LEU LEU
GLY GLY
LYS LYS
ASN ASN
GLU GLU
GLY GLY
ALA ALA
PRO PRO
GLN GLN
GLY GLY
ILE ILE
LEU LEU
GLU GLU
ASP ASP

MET PRO
VAL VAL
ASP ASP
PRO PRO
ASP ASP
ASN ASN
GLU GLU
TYR TYR
LYS LYS
ALA ALA
MET MET
PRO PRO
SER SER
GLU GLU
GLY GLY
TYR TYR
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
ALA ALA

● Molecule 1: Alpha-synuclein



MET ASP
PRO VAL
ASP PHE
PRO MET
ASP LYS
ASN GLY
GLU LEU
TYR SER
LYS TYR
ALA ALA
MET LYS
GLU PRO
SER GLU
G14 G25
VAL VAL
ALA ALA
GLU GLU
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
GLU GLU
G36 E57 K58 Q62 V62 A91
THR THR
GLY GLY
PHE PHE
VAL VAL
LYS LYS
LYS LYS
ASP ASP
GLN GLN
LEU LEU
GLY GLY
LYS LYS
ASN ASN
GLU GLU
GLY GLY
ALA ALA
PRO PRO
GLN GLN
GLY GLY
ILE ILE
LEU LEU
GLU GLU
ASP ASP

MET PRO
VAL VAL
ASP ASP
PRO PRO
ASP ASP
ASN ASN
GLU GLU
TYR TYR
LYS LYS
ALA ALA
MET MET
PRO PRO
SER SER
GLU GLU
GLY GLY
TYR TYR
GLN GLN
ASP ASP
TYR TYR
ALA ALA
GLY GLY
LYS LYS
THR THR
LYS LYS
ALA ALA

4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-0.86°, rise=4.78 Å, axial sym=C2	Depositor
Number of segments used	33479	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	32.6	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.066	Depositor
Minimum map value	-0.045	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	281.6, 281.6, 281.6	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	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	0.69	0/458	1.09	0/618
1	B	0.72	0/458	0.99	0/618
1	C	0.72	0/458	0.98	0/618
1	D	0.70	0/458	0.94	0/618
1	E	0.73	0/458	0.98	0/618
1	F	0.81	0/458	1.06	0/618
1	G	0.70	0/458	1.09	0/618
1	H	0.72	0/458	1.04	0/618
1	I	0.68	0/458	0.97	0/618
1	J	0.71	0/458	0.97	0/618
1	K	0.73	0/458	0.98	0/618
1	L	0.83	0/458	1.08	0/618
All	All	0.73	0/5496	1.02	0/7416

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 [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	458	483	481	0	0
1	B	458	483	481	0	0
1	C	458	483	481	0	0
1	D	458	483	481	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	458	483	481	0	0
1	F	458	483	481	0	0
1	G	458	483	481	0	0
1	H	458	483	481	0	0
1	I	458	483	481	0	0
1	J	458	483	481	0	0
1	K	458	483	481	0	0
1	L	458	483	481	0	0
All	All	5496	5796	5772	0	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 0.

There are no clashes within the asymmetric unit.

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	64/140 (46%)	59 (92%)	5 (8%)	0	100	100
1	B	64/140 (46%)	59 (92%)	5 (8%)	0	100	100
1	C	64/140 (46%)	57 (89%)	7 (11%)	0	100	100
1	D	64/140 (46%)	59 (92%)	5 (8%)	0	100	100
1	E	64/140 (46%)	58 (91%)	6 (9%)	0	100	100
1	F	64/140 (46%)	56 (88%)	8 (12%)	0	100	100
1	G	64/140 (46%)	59 (92%)	5 (8%)	0	100	100
1	H	64/140 (46%)	59 (92%)	5 (8%)	0	100	100
1	I	64/140 (46%)	58 (91%)	6 (9%)	0	100	100
1	J	64/140 (46%)	57 (89%)	7 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	64/140 (46%)	58 (91%)	6 (9%)	0	100	100
1	L	64/140 (46%)	57 (89%)	7 (11%)	0	100	100
All	All	768/1680 (46%)	696 (91%)	72 (9%)	0	100	100

There are no Ramachandran outliers to report.

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	45/103 (44%)	45 (100%)	0	100	100
1	B	45/103 (44%)	45 (100%)	0	100	100
1	C	45/103 (44%)	45 (100%)	0	100	100
1	D	45/103 (44%)	44 (98%)	1 (2%)	45	64
1	E	45/103 (44%)	45 (100%)	0	100	100
1	F	45/103 (44%)	45 (100%)	0	100	100
1	G	45/103 (44%)	45 (100%)	0	100	100
1	H	45/103 (44%)	45 (100%)	0	100	100
1	I	45/103 (44%)	45 (100%)	0	100	100
1	J	45/103 (44%)	45 (100%)	0	100	100
1	K	45/103 (44%)	45 (100%)	0	100	100
1	L	45/103 (44%)	45 (100%)	0	100	100
All	All	540/1236 (44%)	539 (100%)	1 (0%)	85	87

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	63	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	B	65	ASN
1	C	65	ASN
1	D	65	ASN
1	E	65	ASN
1	G	65	ASN
1	H	65	ASN
1	I	65	ASN
1	J	65	ASN
1	K	65	ASN

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.

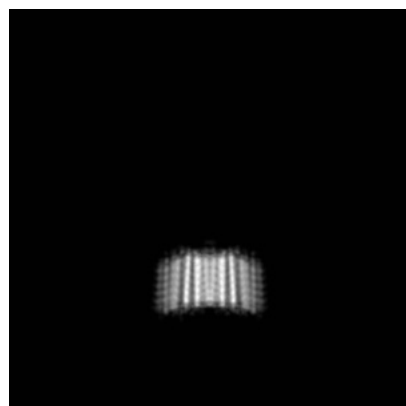
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12266. These allow visual inspection of the internal detail of the map and identification of artifacts.

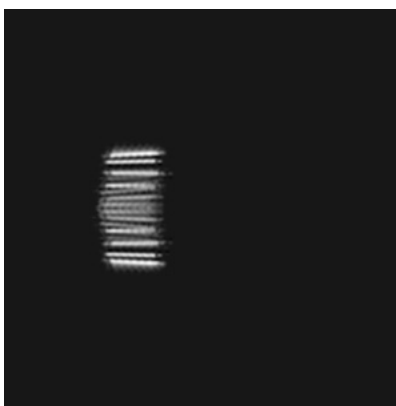
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

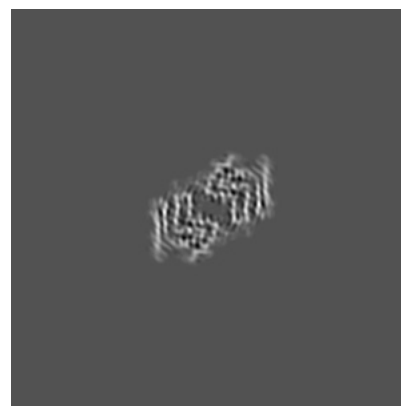
6.1.1 Primary map



X

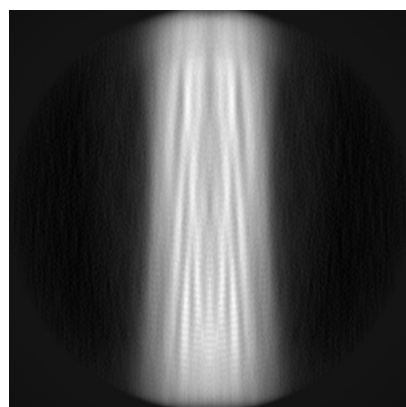


Y

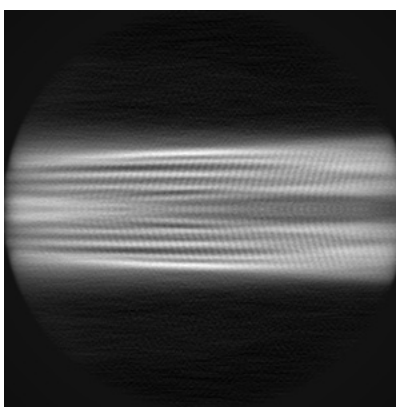


Z

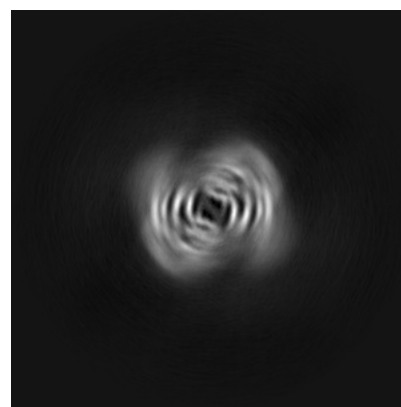
6.1.2 Raw map



X



Y

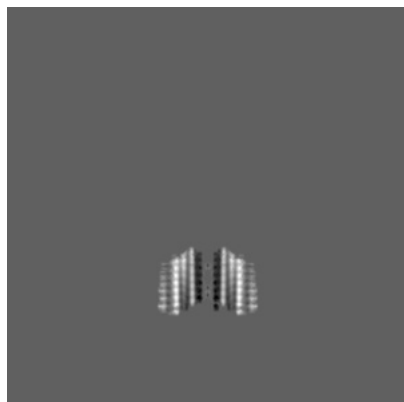


Z

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

6.2 Central slices [i](#)

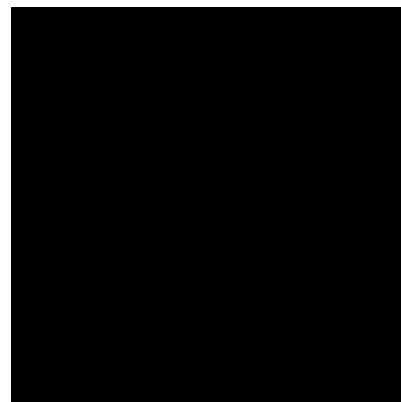
6.2.1 Primary map



X Index: 128

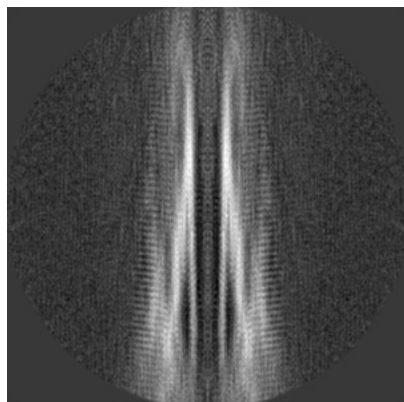


Y Index: 128

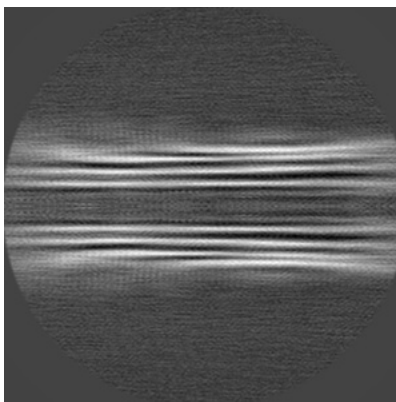


Z Index: 128

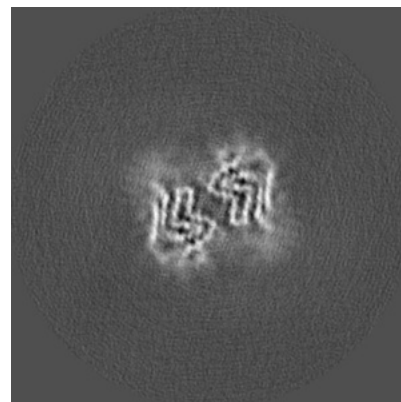
6.2.2 Raw map



X Index: 128



Y Index: 128

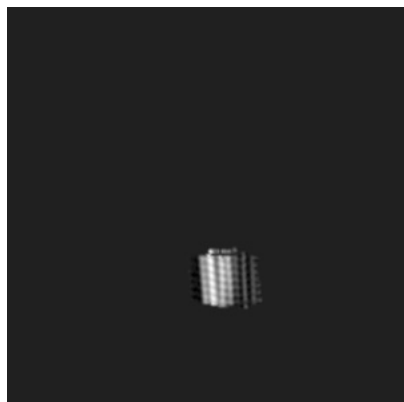


Z Index: 128

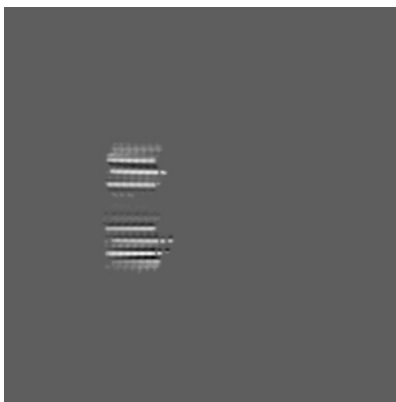
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 158

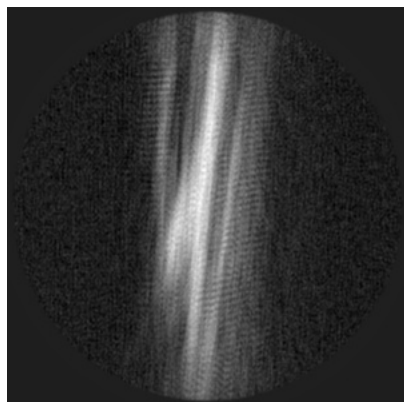


Y Index: 126

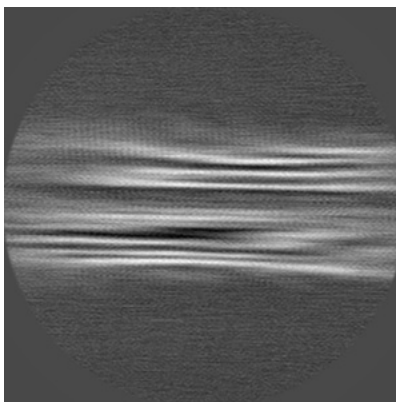


Z Index: 84

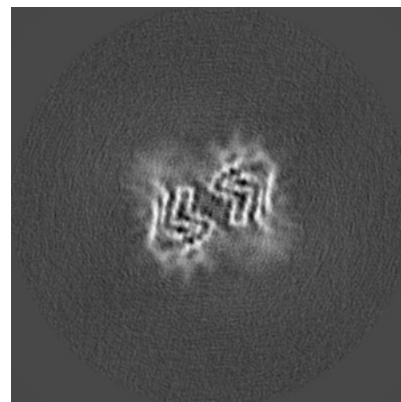
6.3.2 Raw map



X Index: 92



Y Index: 120

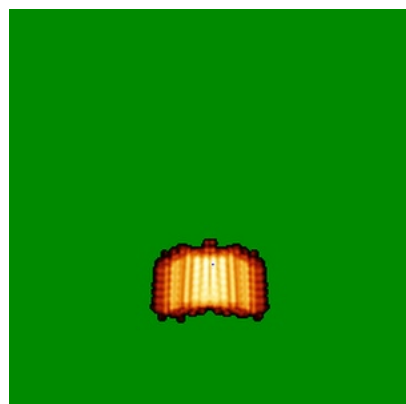


Z Index: 140

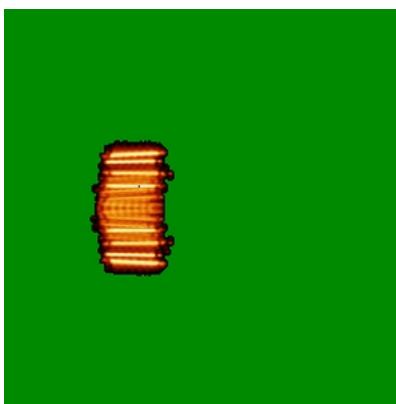
The images above show the largest variance slices of the map in three orthogonal directions.

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

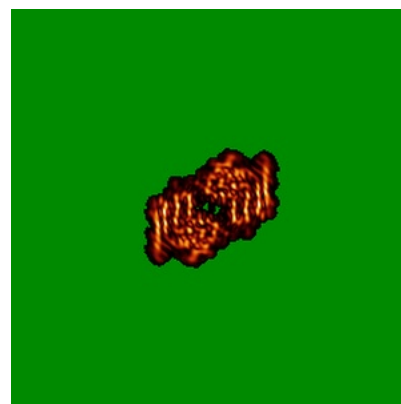
6.4.1 Primary map



X

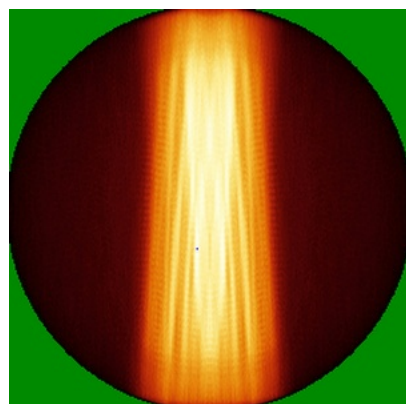


Y

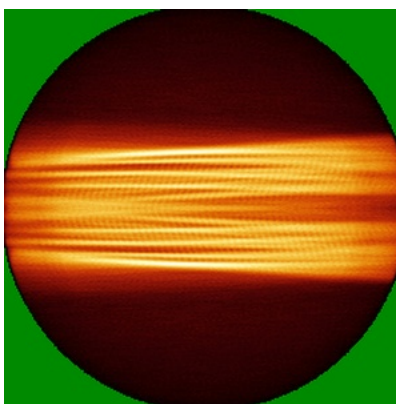


Z

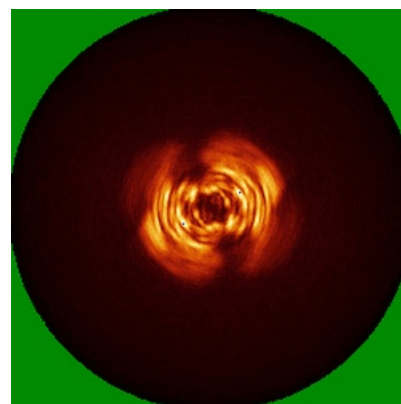
6.4.2 Raw map



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.

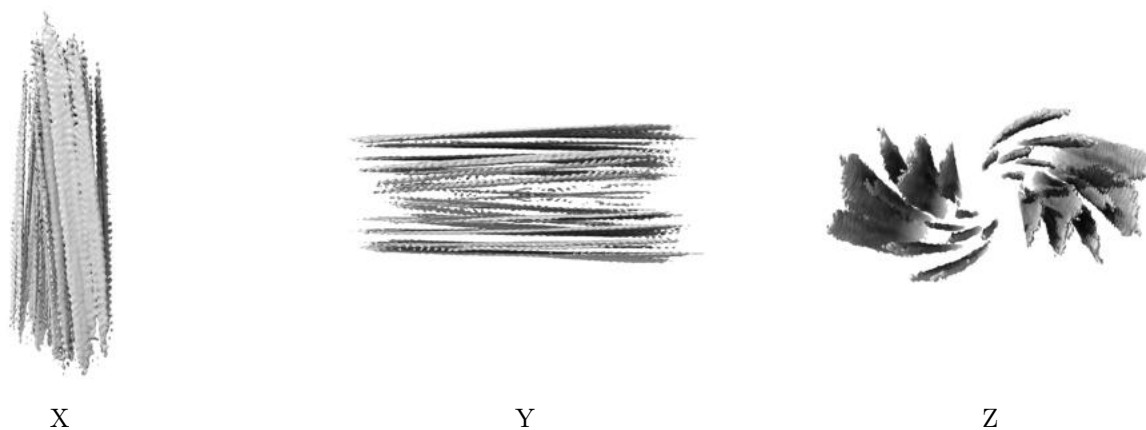
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.01. 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



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.

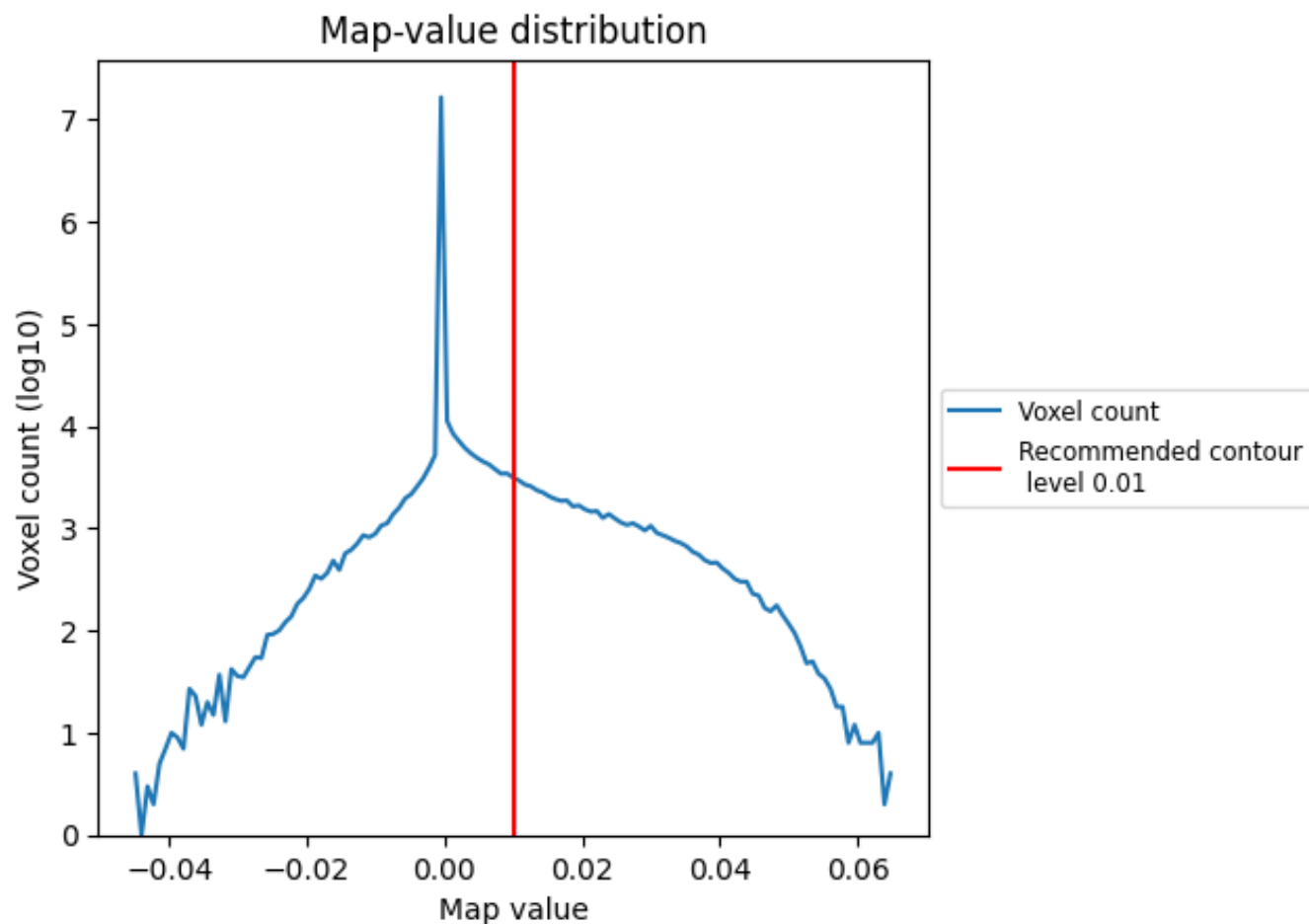
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

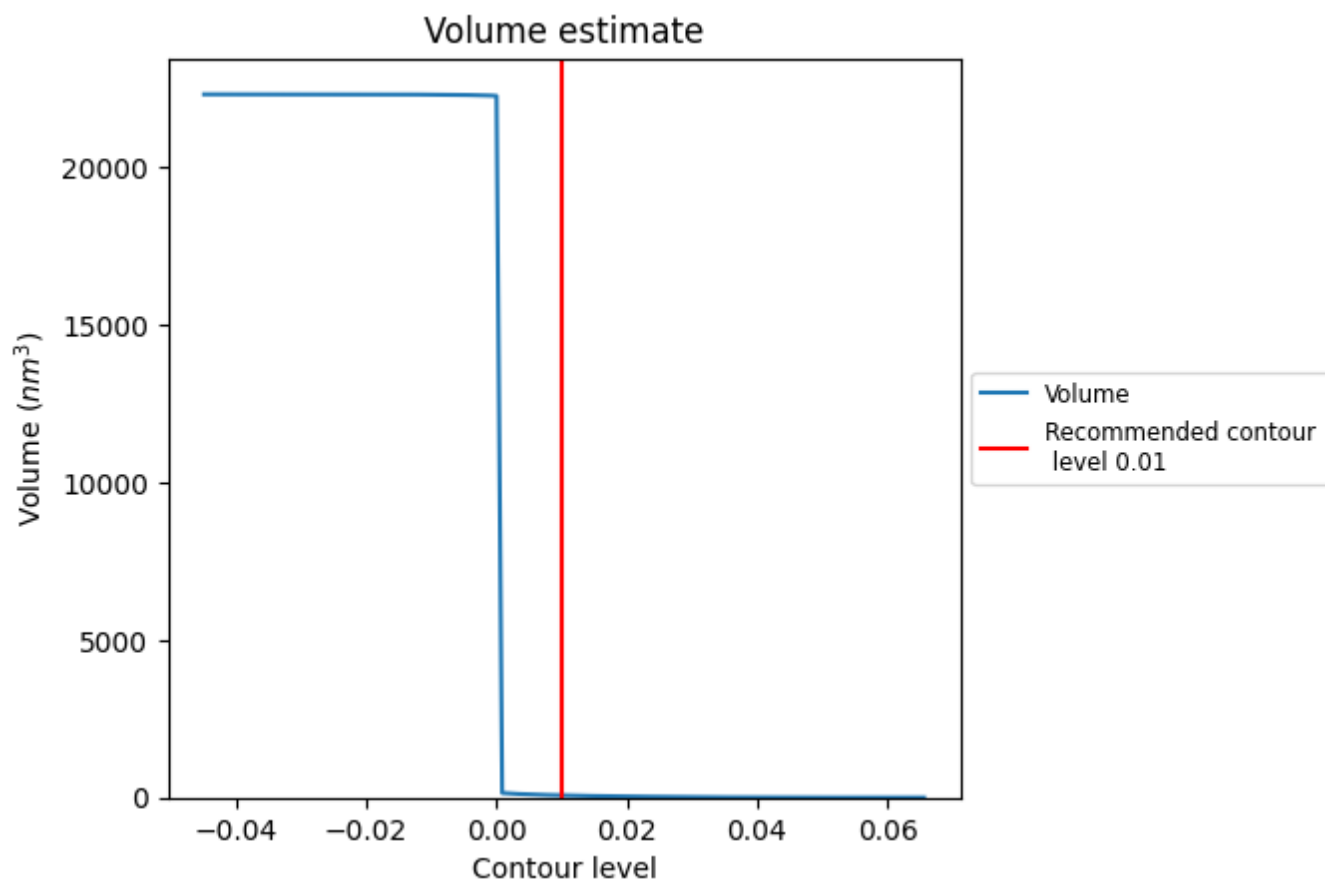
This section contains the results of statistical analysis of the map.

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.

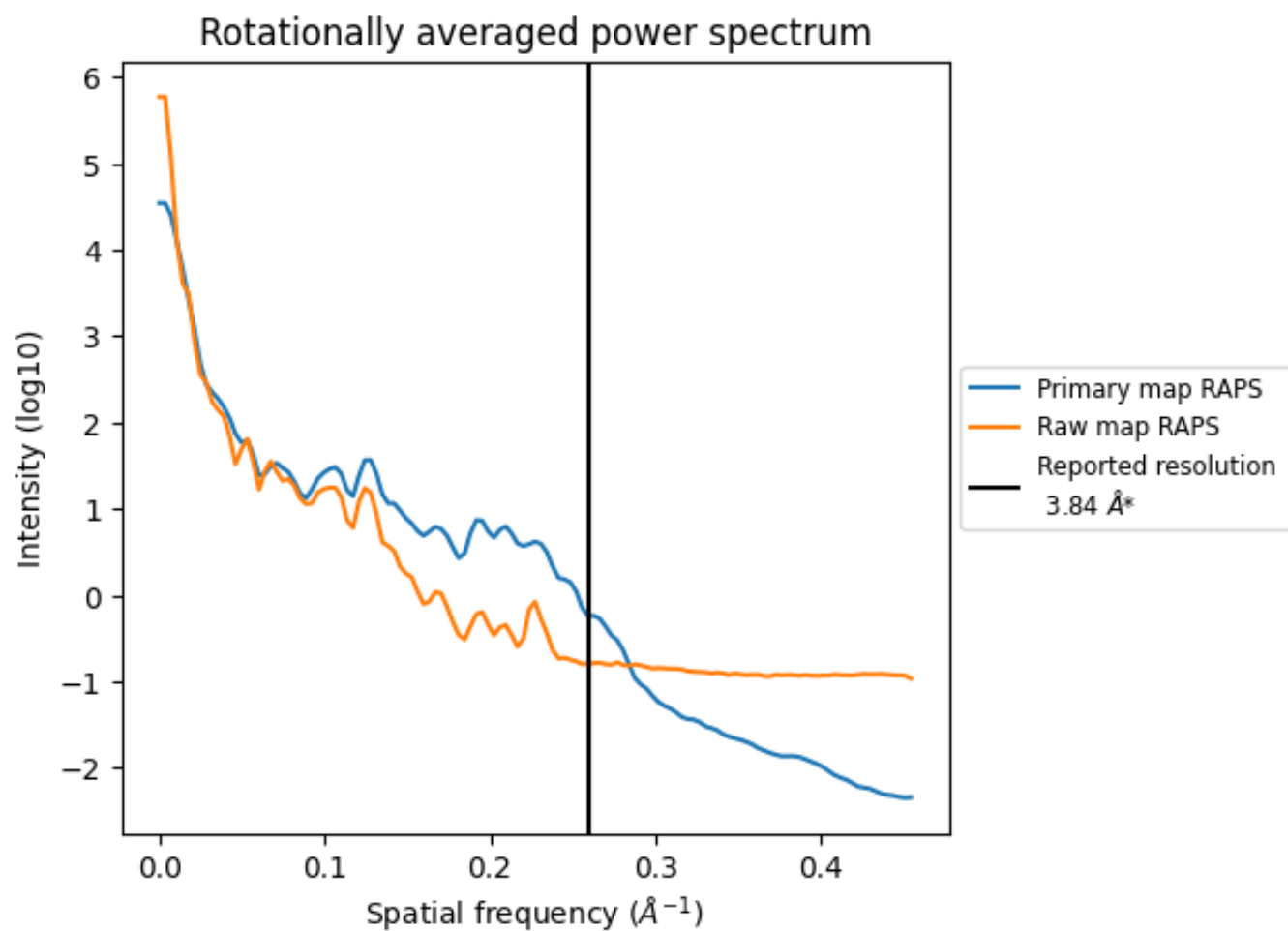
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 70 nm³; this corresponds to an approximate mass of 63 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 ⓘ

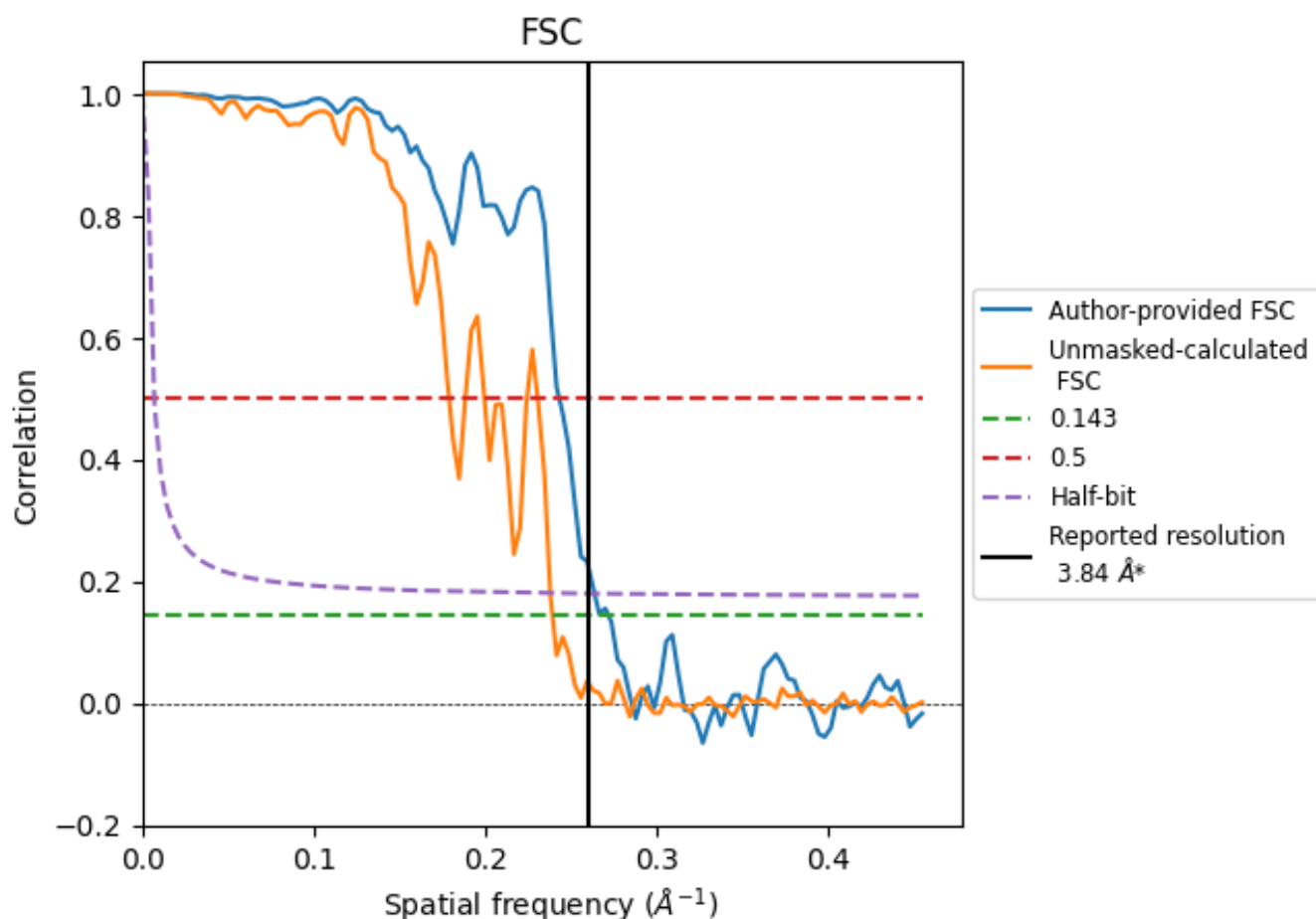


*Reported resolution corresponds to spatial frequency of 0.260 Å⁻¹

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 0.260 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.84	-	-
Author-provided FSC curve	3.68	4.12	3.79
Unmasked-calculated*	4.18	5.59	4.21

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

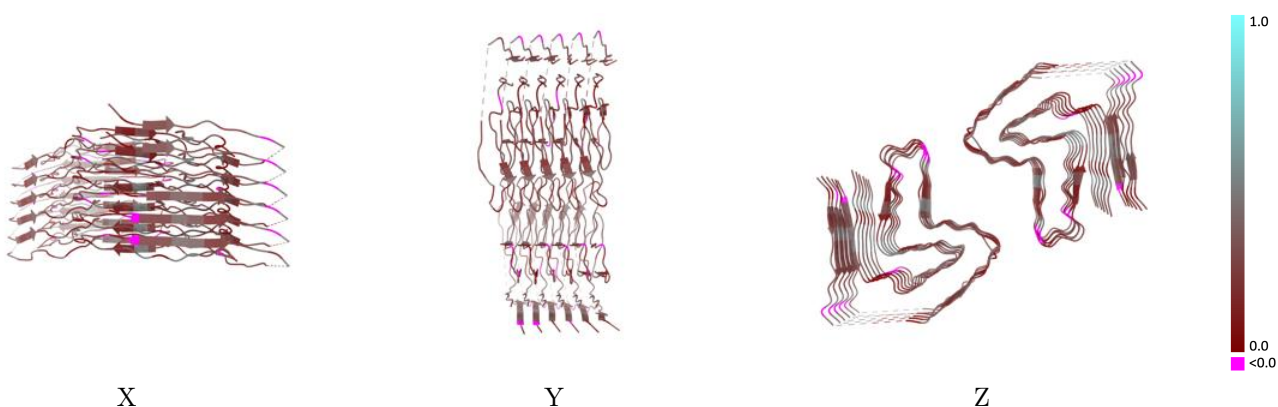
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-12266 and PDB model 7NCH. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)

This section was not generated.

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

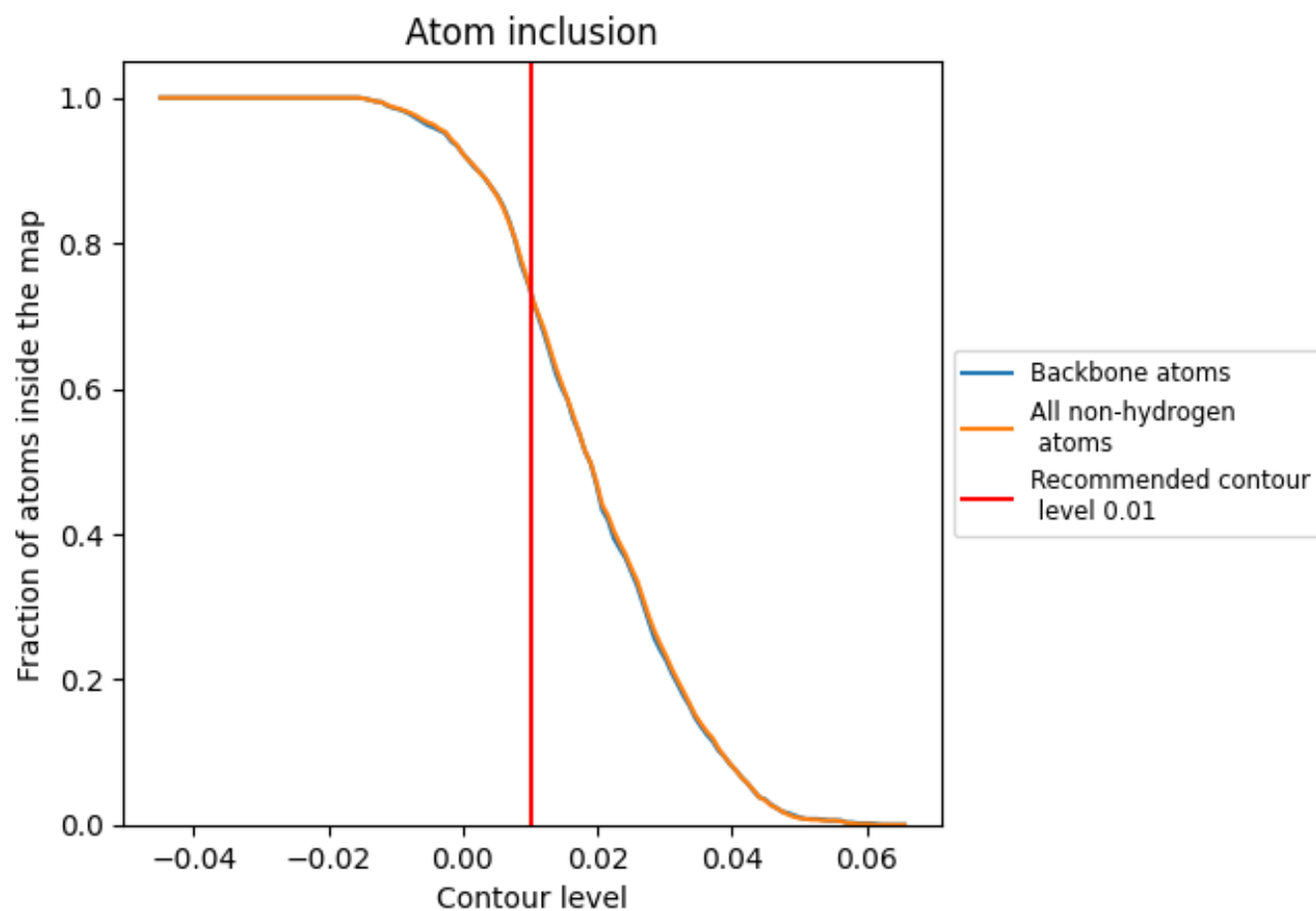


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 74% 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	0.7380	0.2440
A	0.7440	0.2480
B	0.7460	0.2410
C	0.7460	0.2310
D	0.7590	0.2490
E	0.7400	0.2380
F	0.7530	0.2560
G	0.7550	0.2440
H	0.7480	0.2450
I	0.7460	0.2350
J	0.7570	0.2420
K	0.7330	0.2360
L	0.7530	0.2580

