



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 04:01 PM UTC

PDB ID : 8TVG / pdb_00008tv
EMDB ID : EMD-41641
Title : Langya henipavirus postfusion F protein in complex with the 4G5 Fab, local refinement of the viral membrane distal region
Authors : Wang, Z.; Seattle Structural Genomics Center for Infectious Disease (SSG-CID); Veesler, D.
Deposited on : 2023-08-18
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
Mogul : 2022.3.0, CSD as543be (2022)
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

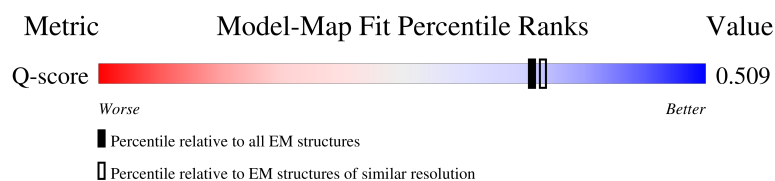
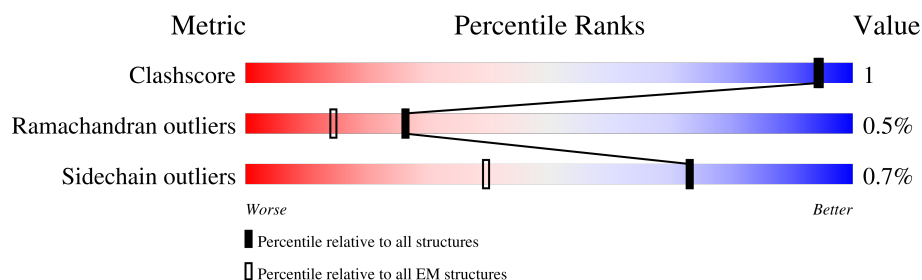
1 Overall quality at a glance

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 $\leq 5\%$. 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	534	9% 66% 29%
1	B	534	9% 66% 29%
1	C	534	9% 66% 29%

2 Entry composition [i](#)

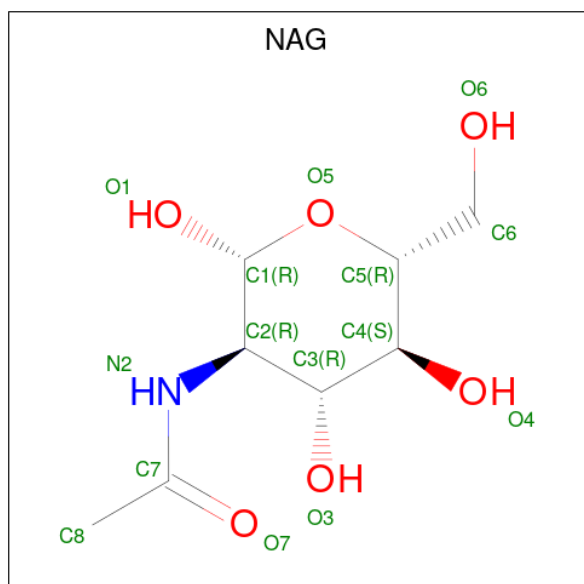
There are 2 unique types of molecules in this entry. The entry contains 8310 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Langya henipavirus postfusion F protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	377	Total	C	N	O	S	0	0
			2742	1763	468	495	16		
1	B	377	Total	C	N	O	S	0	0
			2742	1763	468	495	16		
1	C	377	Total	C	N	O	S	0	0
			2742	1763	468	495	16		

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
2	A	1	Total	C	N	O	0
			14	8	1	5	
2	A	1	Total	C	N	O	0
			14	8	1	5	
2	B	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
2	B	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	

GLY
SER
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● Molecule 1: Langya henipavirus postfusion F protein



MET
ALA
PHE
LEU
THR
LYS
SER
ALA
ILE
ILE
CYS
TYR
LEU
LEU
PHE
TYR
PRO
HIS
ILE
VAL
LYS
SER
L23
K42
S46
I57
D61
G62
V63
R64
M65
E83
P84
D94
K97
S98
GLY
ASN
ASN
LYS
TYR
ARG
PHE
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ALA
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VAL
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LEU
LEU
A159
N160
K161
Q162
G210
N215
D247
E276
F277
V298
D299
G300

D301
D322
R325
E330
L341
C350
D364
T355
S356
K357
C358
A359
R360
V368
G376
D391
T392
D393
T394
S397
Q398
S399
L400
G401
T402
D408
N409
K410
K411
C412
L413
D436
N437
V438
E439
L440
G441
P442
D447
A456
G457
I458
N459
GLN
THR
LEU

GLN
ASN
ALA
GLU
ASP
TYR
ILE
GLU
LYS
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	60440	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	63	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	1700	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.989	Depositor
Minimum map value	-1.678	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.039	Depositor
Recommended contour level	0.5	Depositor
Map size ($\text{\AA}$)	472.08002, 472.08002, 472.08002	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.98350006, 0.98350006, 0.98350006	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	0.66	0/2783	1.03	23/3798 (0.6%)
1	B	0.66	0/2783	1.02	23/3798 (0.6%)
1	C	0.66	0/2783	1.02	23/3798 (0.6%)
All	All	0.66	0/8349	1.03	69/11394 (0.6%)

There are no bond length outliers.

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392	THR	N-CA-C	-7.37	103.88	114.12
1	B	392	THR	N-CA-C	-7.35	103.91	114.12
1	C	392	THR	N-CA-C	-7.33	103.93	114.12
1	C	442	PRO	CA-C-N	6.75	126.71	120.03
1	C	442	PRO	C-N-CA	6.75	126.71	120.03
1	A	442	PRO	CA-C-N	6.73	126.69	120.03
1	A	442	PRO	C-N-CA	6.73	126.69	120.03
1	B	442	PRO	CA-C-N	6.71	126.68	120.03
1	B	442	PRO	C-N-CA	6.71	126.68	120.03
1	A	46	SER	CA-C-N	6.53	127.06	120.14
1	A	46	SER	C-N-CA	6.53	127.06	120.14
1	B	46	SER	CA-C-N	6.52	127.05	120.14
1	B	46	SER	C-N-CA	6.52	127.05	120.14
1	C	46	SER	CA-C-N	6.50	127.03	120.14
1	C	46	SER	C-N-CA	6.50	127.03	120.14
1	A	57	ILE	CA-C-N	6.09	126.06	120.03
1	A	57	ILE	C-N-CA	6.09	126.06	120.03
1	C	277	PHE	CA-C-N	6.08	125.99	119.85
1	C	277	PHE	C-N-CA	6.08	125.99	119.85
1	A	277	PHE	CA-C-N	6.07	125.98	119.85
1	A	277	PHE	C-N-CA	6.07	125.98	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	ILE	CA-C-N	6.07	126.03	120.03
1	B	57	ILE	C-N-CA	6.07	126.03	120.03
1	C	57	ILE	CA-C-N	6.07	126.03	120.03
1	C	57	ILE	C-N-CA	6.07	126.03	120.03
1	B	277	PHE	CA-C-N	6.02	125.93	119.85
1	B	277	PHE	C-N-CA	6.02	125.93	119.85
1	C	368	VAL	CA-C-N	6.02	125.93	119.85
1	C	368	VAL	C-N-CA	6.02	125.93	119.85
1	A	368	VAL	CA-C-N	6.01	125.92	119.85
1	A	368	VAL	C-N-CA	6.01	125.92	119.85
1	B	368	VAL	CA-C-N	5.99	125.90	119.85
1	B	368	VAL	C-N-CA	5.99	125.90	119.85
1	A	83	GLU	CA-C-N	5.90	125.58	119.56
1	A	83	GLU	C-N-CA	5.90	125.58	119.56
1	C	83	GLU	CA-C-N	5.90	125.58	119.56
1	C	83	GLU	C-N-CA	5.90	125.58	119.56
1	B	83	GLU	CA-C-N	5.89	125.56	119.56
1	B	83	GLU	C-N-CA	5.89	125.56	119.56
1	A	210	GLY	CA-C-N	5.83	125.96	119.32
1	A	210	GLY	C-N-CA	5.83	125.96	119.32
1	C	210	GLY	CA-C-N	5.76	125.89	119.32
1	C	210	GLY	C-N-CA	5.76	125.89	119.32
1	B	441	GLY	CA-C-N	5.75	126.30	120.38
1	B	441	GLY	C-N-CA	5.75	126.30	120.38
1	B	210	GLY	CA-C-N	5.74	125.86	119.32
1	B	210	GLY	C-N-CA	5.74	125.86	119.32
1	A	441	GLY	CA-C-N	5.73	126.28	120.38
1	A	441	GLY	C-N-CA	5.73	126.28	120.38
1	C	341	LEU	CA-C-N	5.72	125.73	119.90
1	C	341	LEU	C-N-CA	5.72	125.73	119.90
1	B	341	LEU	CA-C-N	5.70	125.72	119.90
1	B	341	LEU	C-N-CA	5.70	125.72	119.90
1	C	441	GLY	CA-C-N	5.69	126.24	120.38
1	C	441	GLY	C-N-CA	5.69	126.24	120.38
1	A	341	LEU	CA-C-N	5.68	125.69	119.90
1	A	341	LEU	C-N-CA	5.68	125.69	119.90
1	B	394	THR	CA-C-N	5.56	125.46	119.85
1	B	394	THR	C-N-CA	5.56	125.46	119.85
1	A	394	THR	CA-C-N	5.55	125.46	119.85
1	A	394	THR	C-N-CA	5.55	125.46	119.85
1	C	394	THR	CA-C-N	5.54	125.44	119.85
1	C	394	THR	C-N-CA	5.54	125.44	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	215	ASN	CA-C-N	5.12	125.42	119.47
1	B	215	ASN	C-N-CA	5.12	125.42	119.47
1	C	215	ASN	CA-C-N	5.11	125.40	119.47
1	C	215	ASN	C-N-CA	5.11	125.40	119.47
1	A	215	ASN	CA-C-N	5.11	125.40	119.47
1	A	215	ASN	C-N-CA	5.11	125.40	119.47

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	2742	0	2644	7	0
1	B	2742	0	2644	6	0
1	C	2742	0	2644	6	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
2	C	28	0	26	0	0
All	All	8310	0	8010	19	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 1.

All (19) 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:355:THR:HA	1:C:358:CYS:SG	2.40	0.61
1:A:355:THR:HA	1:A:358:CYS:SG	2.40	0.61
1:B:355:THR:HA	1:B:358:CYS:SG	2.40	0.60
1:B:358:CYS:O	1:B:359:ALA:C	2.47	0.57
1:A:358:CYS:O	1:A:359:ALA:C	2.47	0.57
1:C:358:CYS:O	1:C:359:ALA:C	2.47	0.57
1:A:42:LYS:NZ	1:A:247:ASP:OD1	2.45	0.49
1:B:42:LYS:NZ	1:B:247:ASP:OD1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:LYS:NZ	1:C:247:ASP:OD1	2.45	0.48
1:C:97:LYS:NZ	1:C:276:GLU:OE2	2.48	0.47
1:A:97:LYS:NZ	1:A:276:GLU:OE2	2.48	0.47
1:B:97:LYS:NZ	1:B:276:GLU:OE2	2.48	0.46
1:A:336:ASP:O	1:A:337:ASN:HB2	2.16	0.46
1:B:350:CYS:SG	1:B:358:CYS:N	2.91	0.43
1:C:350:CYS:SG	1:C:358:CYS:N	2.91	0.43
1:B:83:GLU:N	1:B:84:PRO:CD	2.82	0.43
1:A:350:CYS:SG	1:A:358:CYS:N	2.91	0.43
1:A:83:GLU:N	1:A:84:PRO:CD	2.82	0.42
1:C:83:GLU:N	1:C:84:PRO:CD	2.82	0.42

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	373/534 (70%)	361 (97%)	10 (3%)	2 (0%)	24	55
1	B	373/534 (70%)	361 (97%)	10 (3%)	2 (0%)	24	55
1	C	373/534 (70%)	361 (97%)	10 (3%)	2 (0%)	24	55
All	All	1119/1602 (70%)	1083 (97%)	30 (3%)	6 (0%)	26	55

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	359	ALA
1	B	359	ALA
1	C	359	ALA
1	A	298	VAL
1	B	298	VAL
1	C	298	VAL

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	268/460 (58%)	266 (99%)	2 (1%)	76	80
1	B	268/460 (58%)	266 (99%)	2 (1%)	76	80
1	C	268/460 (58%)	266 (99%)	2 (1%)	76	80
All	All	804/1380 (58%)	798 (99%)	6 (1%)	73	80

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	357	LYS
1	A	459	ASN
1	B	357	LYS
1	B	459	ASN
1	C	357	LYS
1	C	459	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	95	ASN
1	A	279	ASN
1	A	337	ASN
1	B	70	GLN
1	B	95	ASN
1	B	279	ASN
1	C	70	GLN
1	C	279	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](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	B	602	1	14,14,15	1.62	2 (14%)	17,19,21	0.95	1 (5%)
2	NAG	C	601	1	14,14,15	1.67	2 (14%)	17,19,21	0.97	1 (5%)
2	NAG	B	601	1	14,14,15	1.67	2 (14%)	17,19,21	0.96	1 (5%)
2	NAG	C	602	1	14,14,15	1.61	2 (14%)	17,19,21	0.95	1 (5%)
2	NAG	A	601	1	14,14,15	1.68	2 (14%)	17,19,21	0.96	1 (5%)
2	NAG	A	602	1	14,14,15	1.61	2 (14%)	17,19,21	0.94	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	602	1	-	1/6/23/26	0/1/1/1
2	NAG	C	601	1	-	0/6/23/26	0/1/1/1
2	NAG	B	601	1	-	0/6/23/26	0/1/1/1
2	NAG	C	602	1	-	1/6/23/26	0/1/1/1
2	NAG	A	601	1	-	0/6/23/26	0/1/1/1
2	NAG	A	602	1	-	1/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	NAG	C1-C2	4.71	1.58	1.52
2	C	601	NAG	C1-C2	4.71	1.58	1.52
2	A	601	NAG	C1-C2	4.70	1.58	1.52
2	A	602	NAG	C1-C2	4.56	1.58	1.52
2	B	602	NAG	C1-C2	4.56	1.58	1.52
2	C	602	NAG	C1-C2	4.56	1.58	1.52
2	A	601	NAG	O5-C5	2.60	1.48	1.43
2	B	601	NAG	O5-C5	2.58	1.48	1.43
2	C	601	NAG	O5-C5	2.54	1.48	1.43
2	A	602	NAG	O5-C5	2.32	1.47	1.43
2	C	602	NAG	O5-C5	2.31	1.47	1.43
2	B	602	NAG	O5-C5	2.29	1.47	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	NAG	C8-C7-N2	2.24	119.84	116.12
2	A	601	NAG	C8-C7-N2	2.24	119.83	116.12
2	B	601	NAG	C8-C7-N2	2.23	119.81	116.12
2	A	602	NAG	C8-C7-N2	2.22	119.80	116.12
2	B	602	NAG	C8-C7-N2	2.21	119.78	116.12
2	C	602	NAG	C8-C7-N2	2.21	119.78	116.12

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	NAG	O5-C5-C6-O6
2	B	602	NAG	O5-C5-C6-O6
2	C	602	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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.

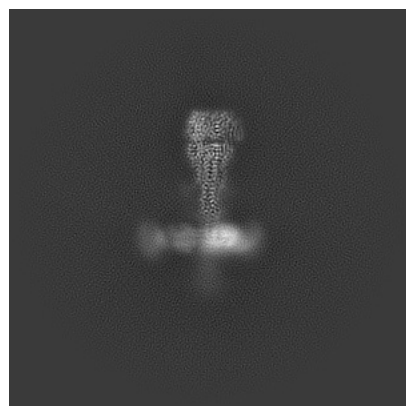
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41641. These allow visual inspection of the internal detail of the map and identification of artifacts.

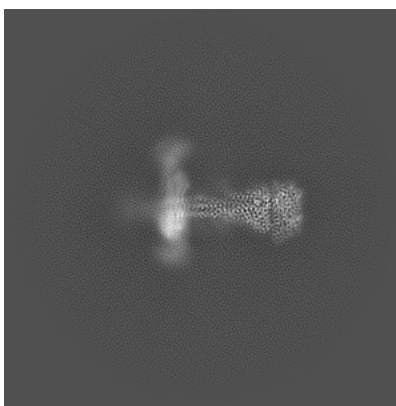
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

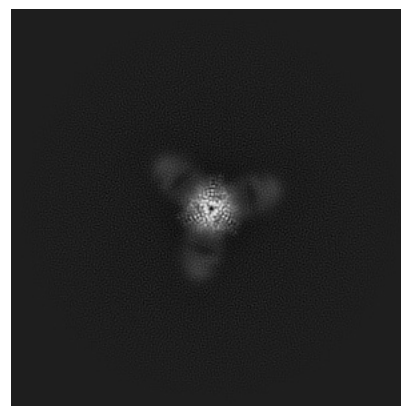
6.1.1 Primary map



X

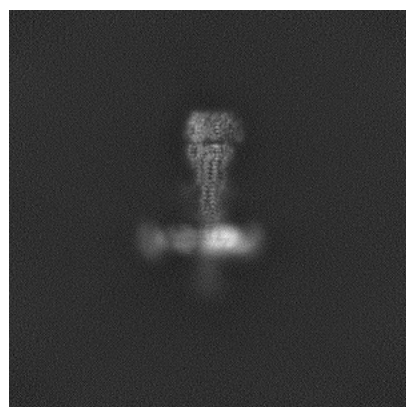


Y

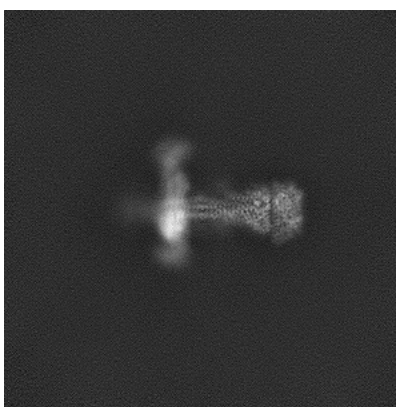


Z

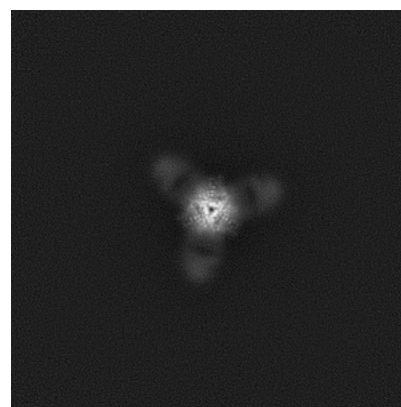
6.1.2 Raw map



X



Y

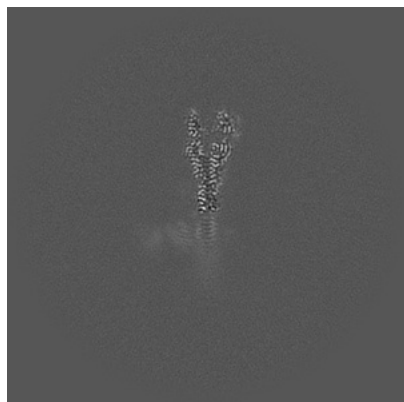


Z

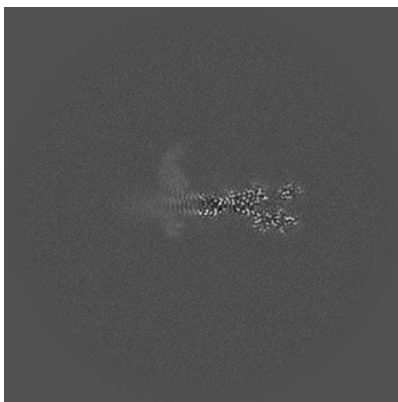
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

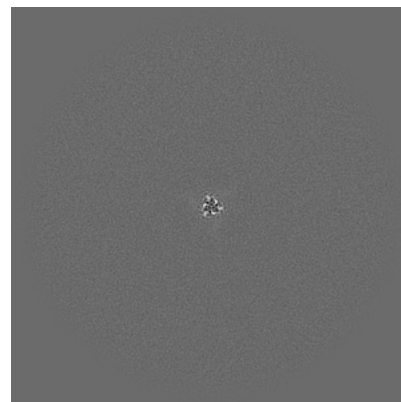
6.2.1 Primary map



X Index: 240



Y Index: 240



Z Index: 240

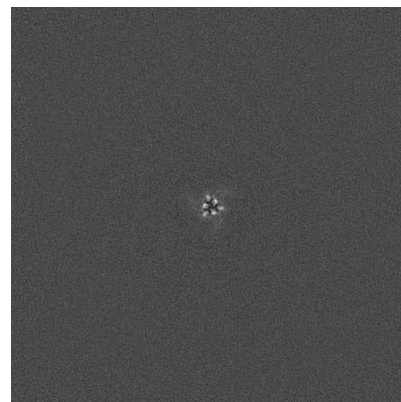
6.2.2 Raw map



X Index: 240



Y Index: 240

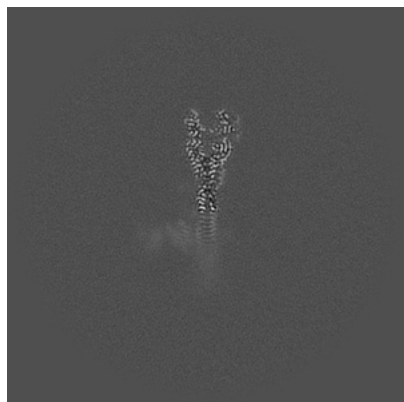


Z Index: 240

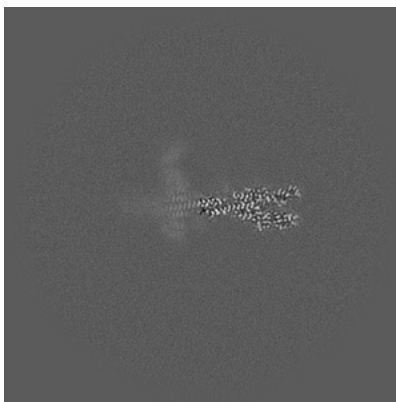
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

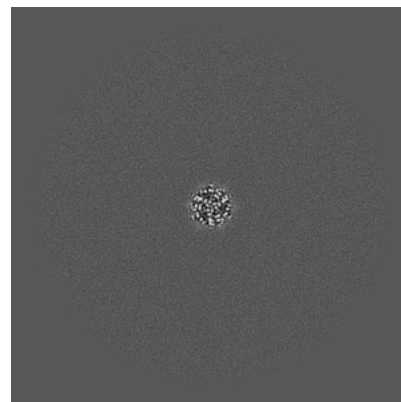
6.3.1 Primary map



X Index: 241



Y Index: 245



Z Index: 302

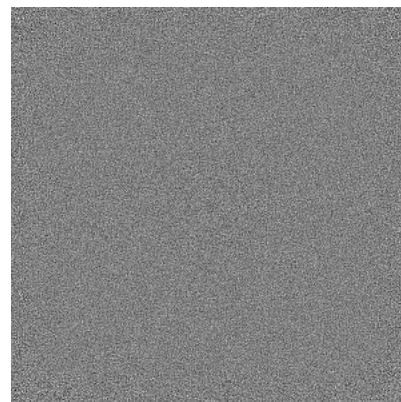
6.3.2 Raw map



X Index: 241



Y Index: 245

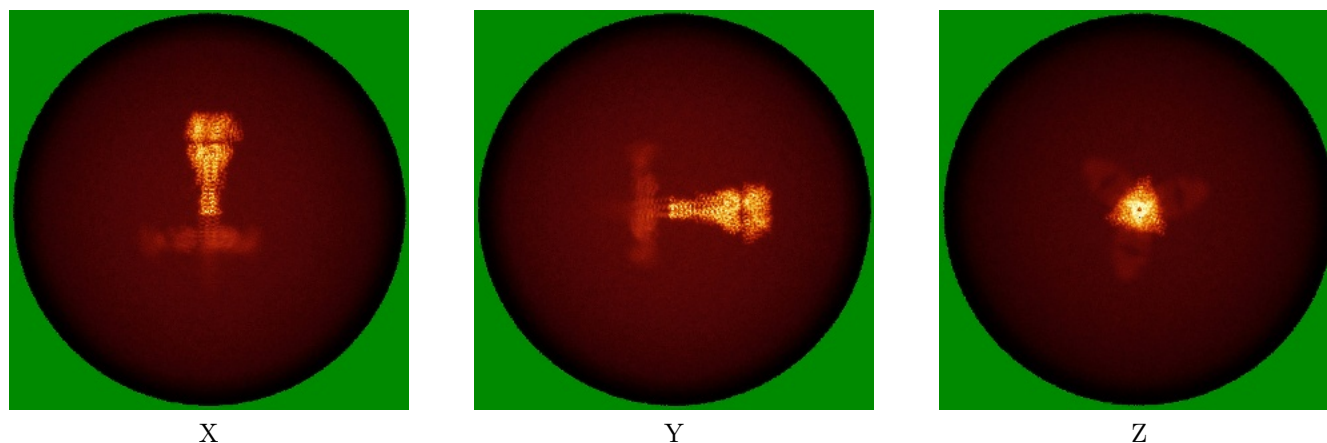


Z Index: 0

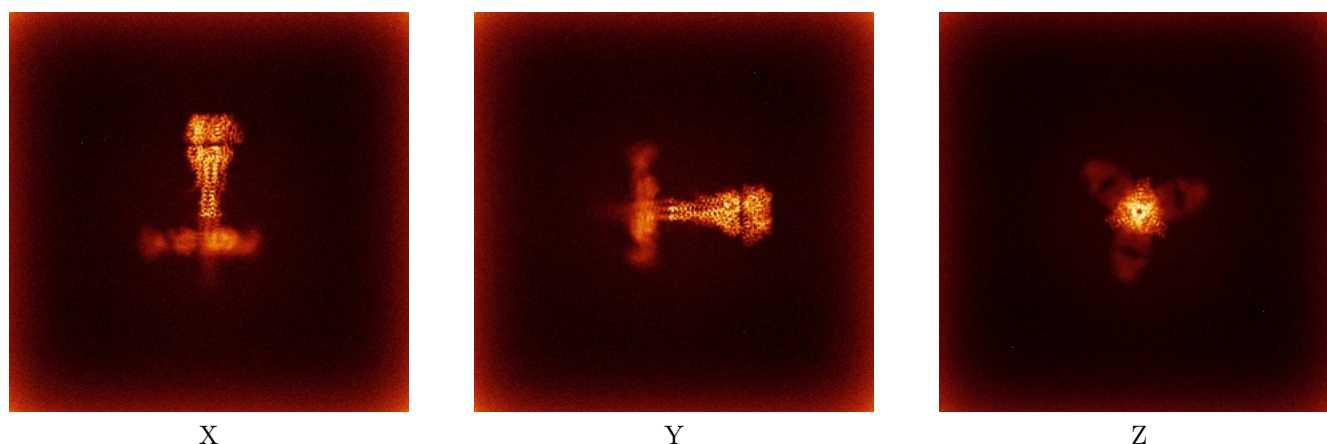
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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.

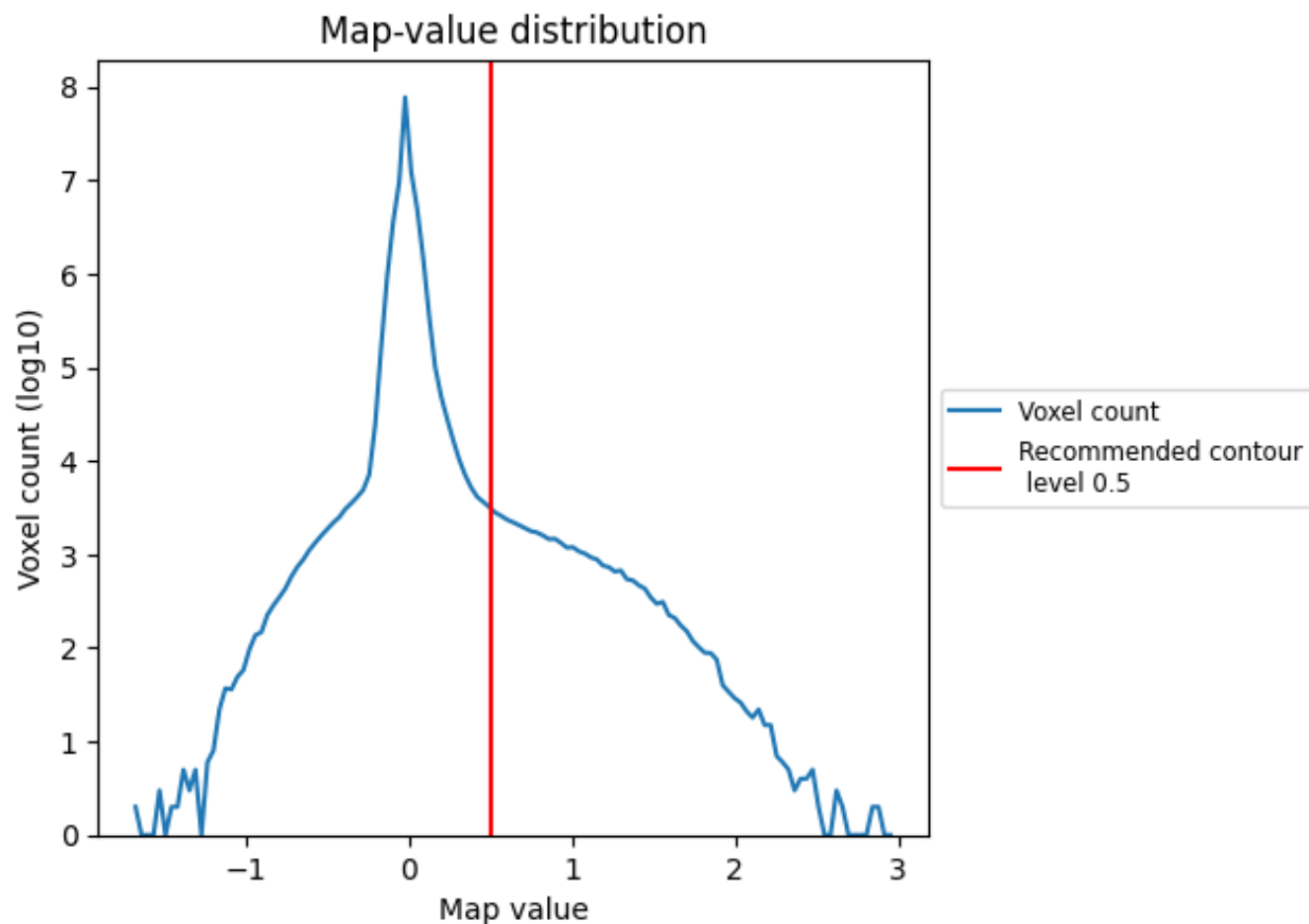
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

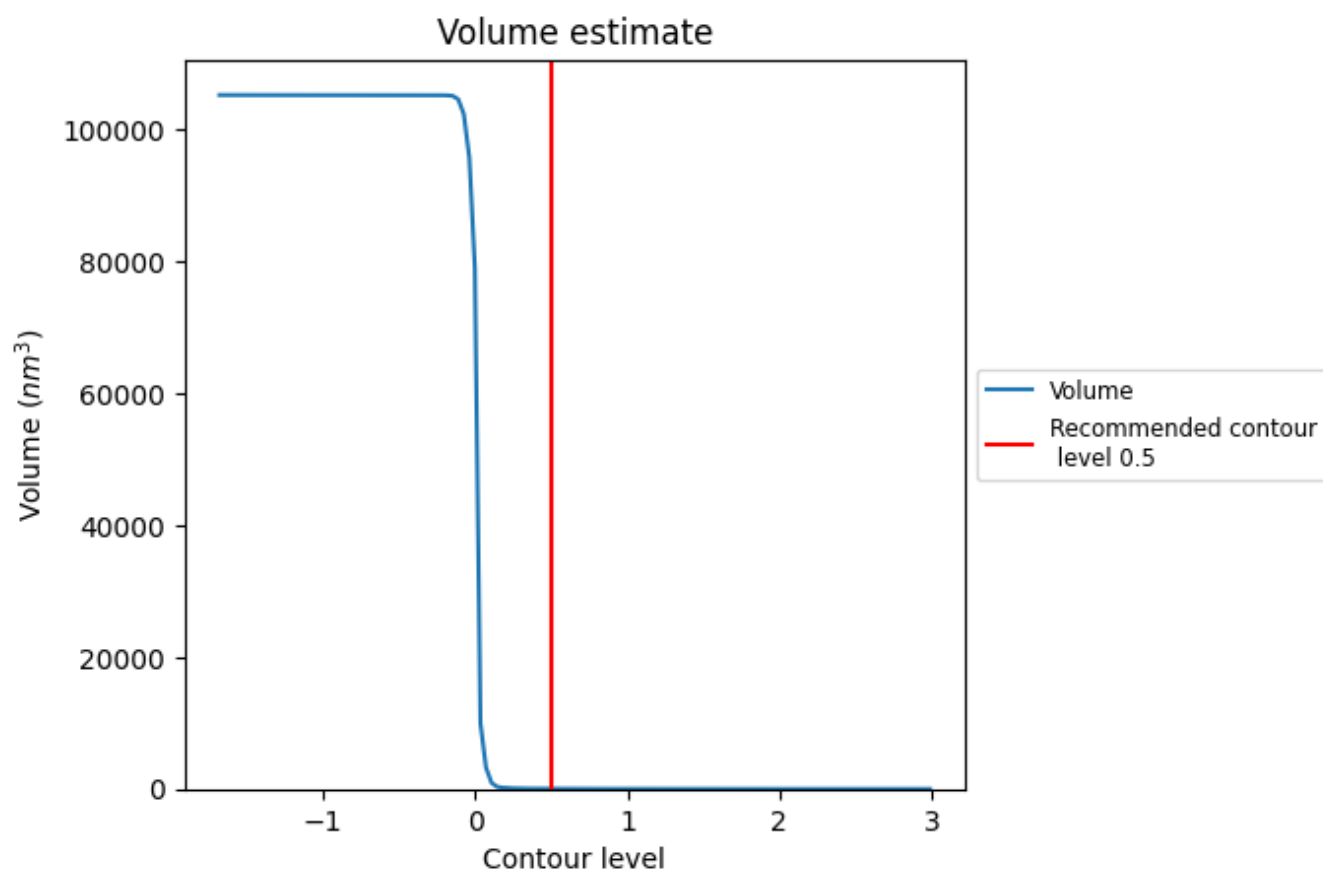
This section contains the results of statistical analysis of the map.

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.

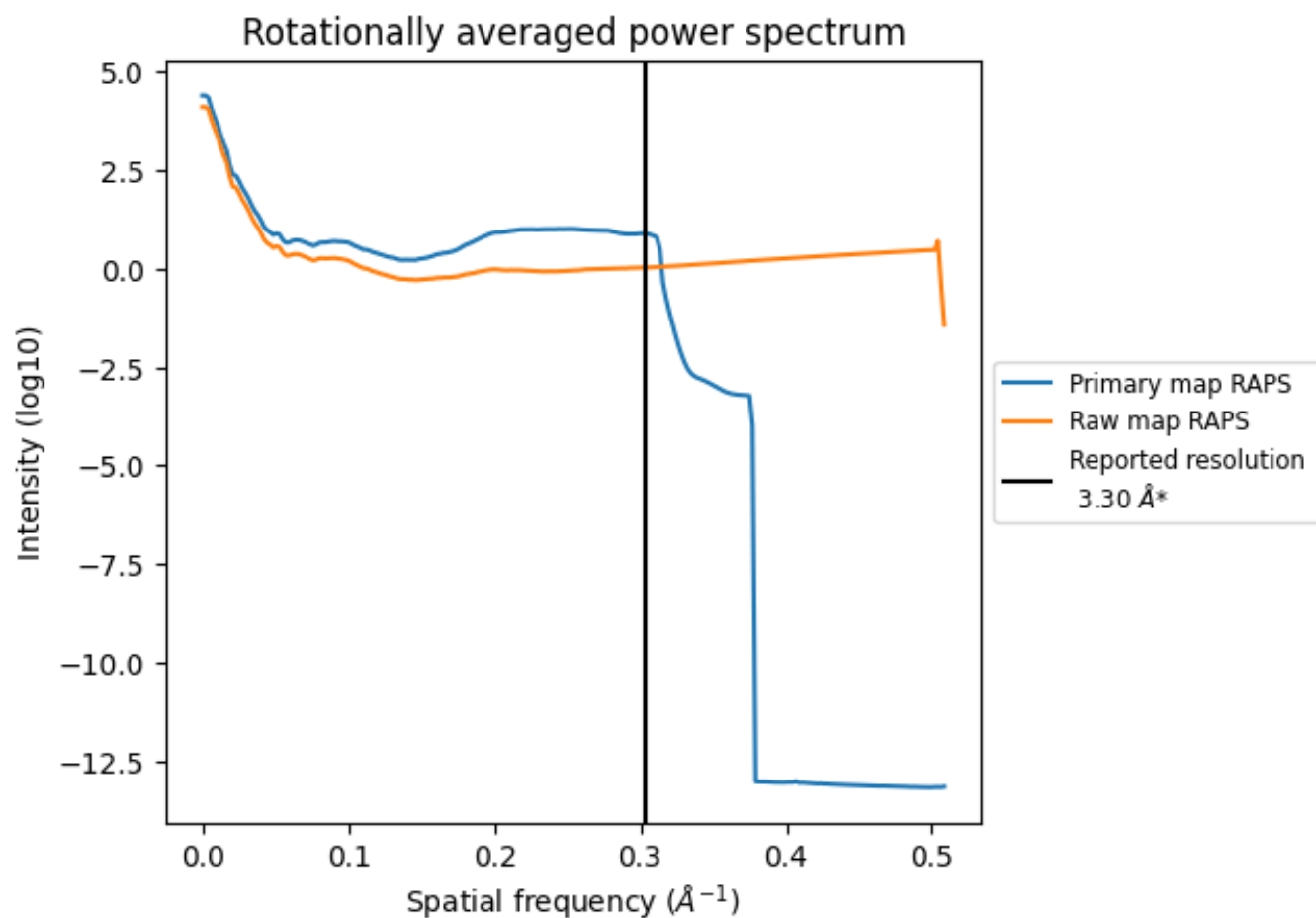
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 37 nm³; this corresponds to an approximate mass of 34 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 ⓘ

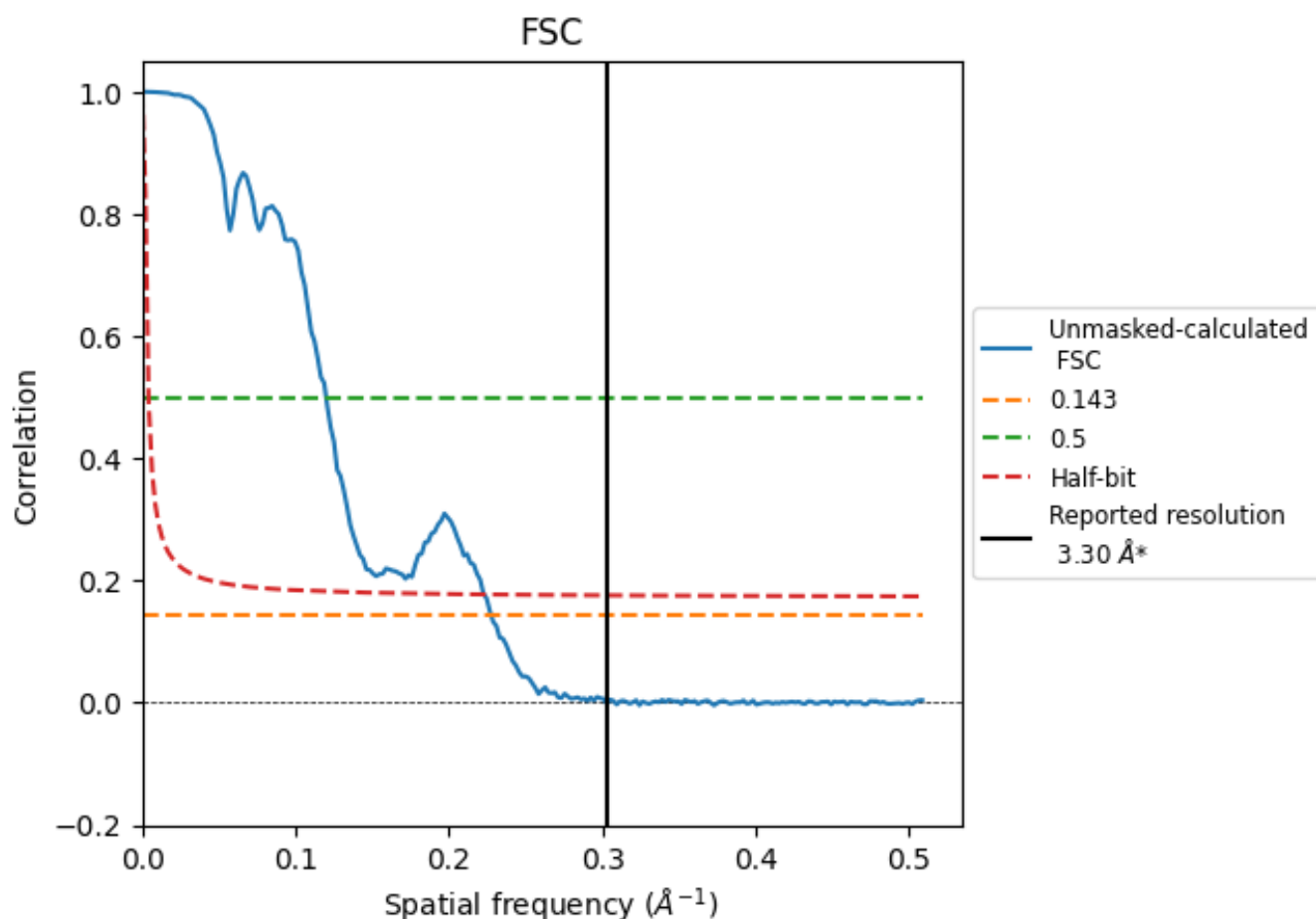


*Reported resolution corresponds to spatial frequency of 0.303 Å⁻¹

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 0.303 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

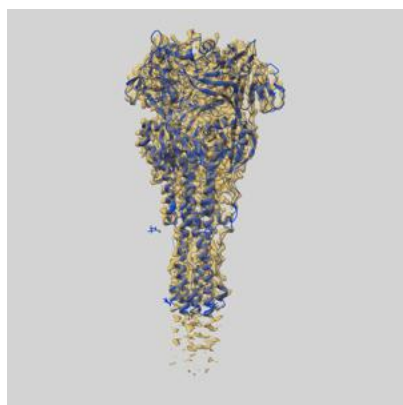
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.40	8.34	4.49

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.40 differs from the reported value 3.3 by more than 10 %

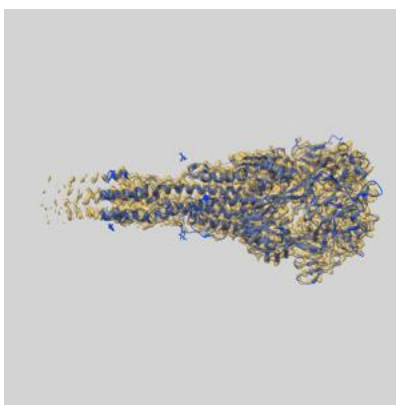
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-41641 and PDB model 8TVG. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

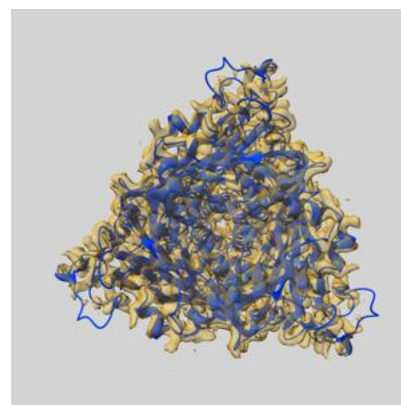
9.1 Map-model overlay [i](#)



X



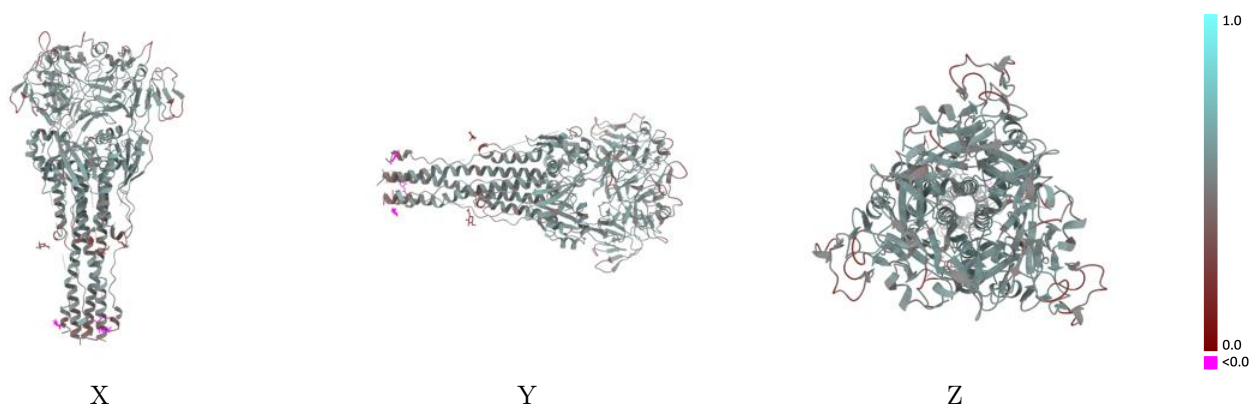
Y



Z

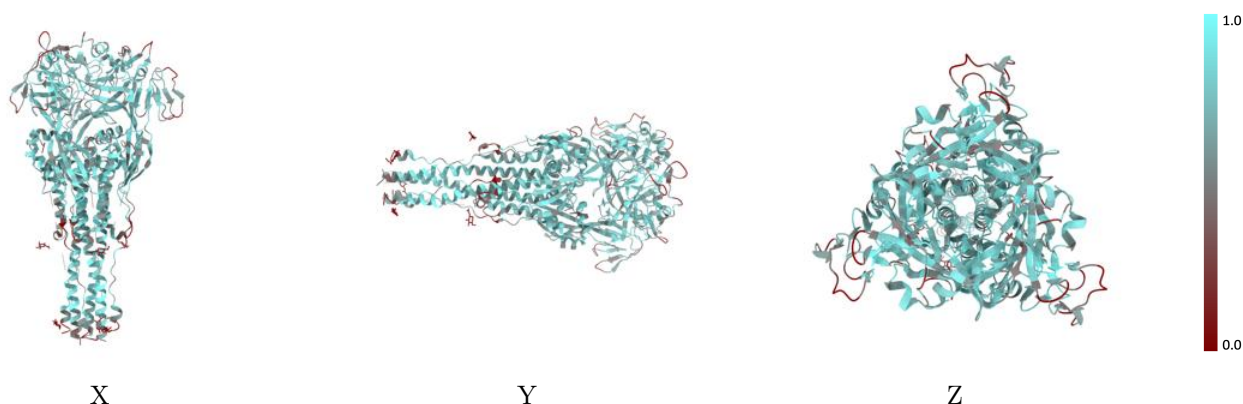
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



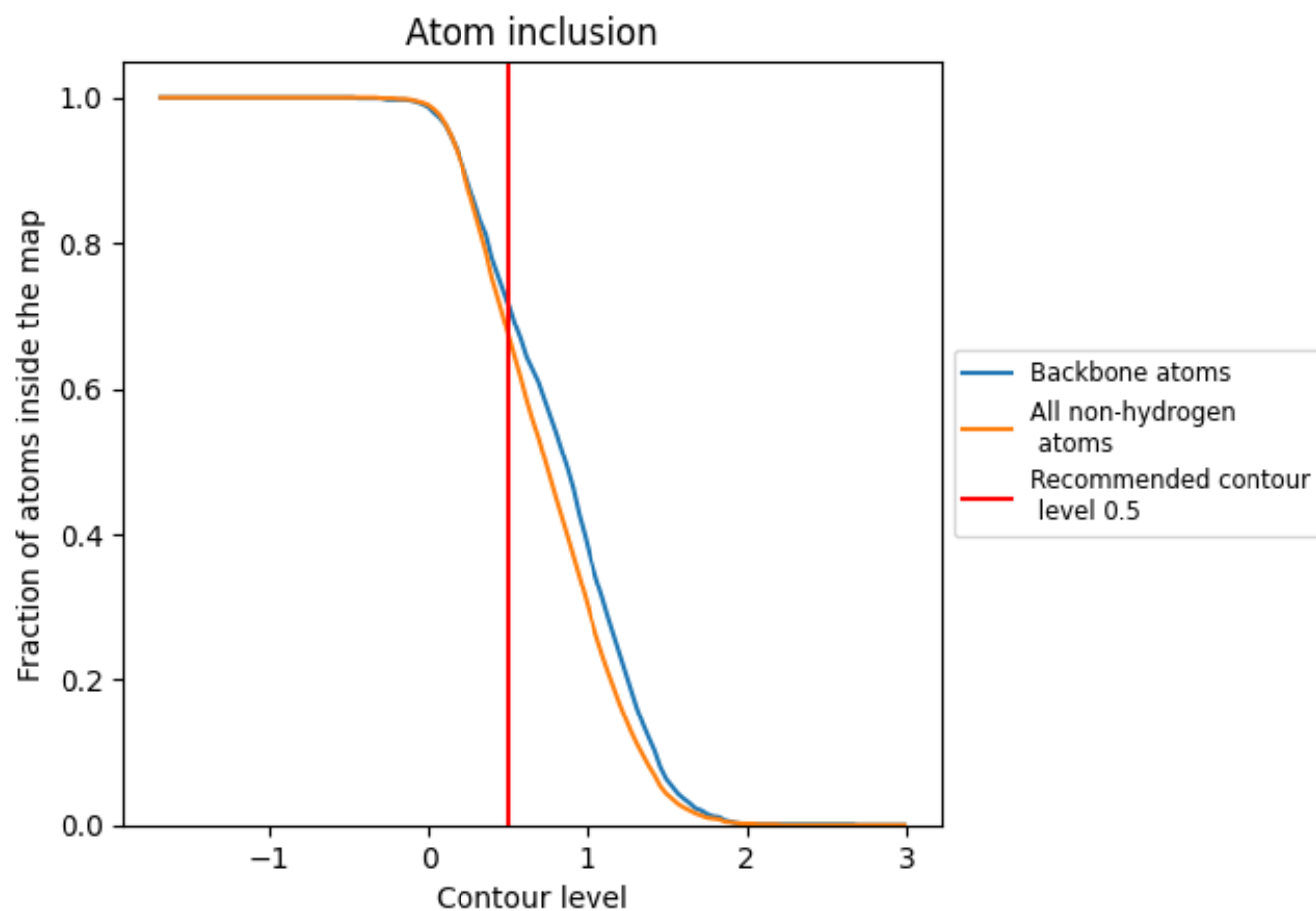
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 72% of all backbone atoms, 68% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6780	0.5090
A	0.6780	0.5090
B	0.6770	0.5100
C	0.6800	0.5100

