



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 10:59 AM UTC

PDB ID : 9MHA / pdb_00009mha
EMDB ID : EMD-48271
Title : CryoEM Structure of Zaire Ebola Virus Envelope Glycoprotein GP
Authors : Weidle, C.; Borst, A.J.
Deposited on : 2024-12-11
Resolution : 3.05 Å(reported)
Based on initial model : 5JQ3

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:

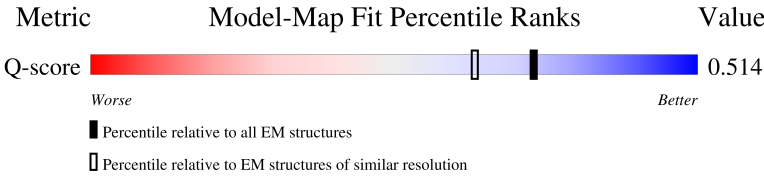
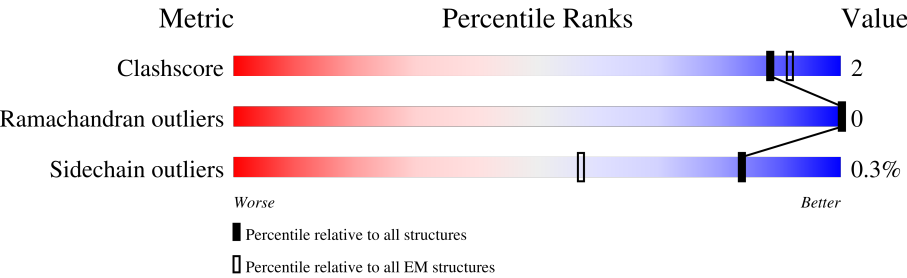
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.05 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.



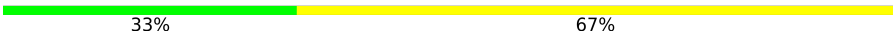
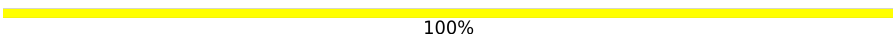
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13971 (2.55 - 3.55)

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	C	168	
1	E	168	
1	F	168	
2	A	321	

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Mol	Chain	Length	Quality of chain
2	B	321	 53% 45%
2	D	321	 51% 45%
3	G	3	 33% 67%
3	H	3	 33% 67%
3	I	3	 100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6188 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 GP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	95	Total	C	N	O	S	0	0
			749	483	132	131	3		
1	E	95	Total	C	N	O	S	0	0
			749	483	131	132	3		
1	F	95	Total	C	N	O	S	0	0
			752	484	132	133	3		

There are 111 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	633	GLY	-	expression tag	UNP Q05320
C	634	SER	-	expression tag	UNP Q05320
C	635	GLY	-	expression tag	UNP Q05320
C	636	TYR	-	expression tag	UNP Q05320
C	637	ILE	-	expression tag	UNP Q05320
C	638	PRO	-	expression tag	UNP Q05320
C	639	GLU	-	expression tag	UNP Q05320
C	640	ALA	-	expression tag	UNP Q05320
C	641	PRO	-	expression tag	UNP Q05320
C	642	ARG	-	expression tag	UNP Q05320
C	643	ASP	-	expression tag	UNP Q05320
C	644	GLY	-	expression tag	UNP Q05320
C	645	GLN	-	expression tag	UNP Q05320
C	646	ALA	-	expression tag	UNP Q05320
C	647	TYR	-	expression tag	UNP Q05320
C	648	VAL	-	expression tag	UNP Q05320
C	649	ARG	-	expression tag	UNP Q05320
C	650	LYS	-	expression tag	UNP Q05320
C	651	ASP	-	expression tag	UNP Q05320
C	652	GLY	-	expression tag	UNP Q05320
C	653	GLU	-	expression tag	UNP Q05320
C	654	TRP	-	expression tag	UNP Q05320
C	655	VAL	-	expression tag	UNP Q05320
C	656	LEU	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
C	657	LEU	-	expression tag	UNP Q05320
C	658	SER	-	expression tag	UNP Q05320
C	659	THR	-	expression tag	UNP Q05320
C	660	PHE	-	expression tag	UNP Q05320
C	661	LEU	-	expression tag	UNP Q05320
C	662	GLY	-	expression tag	UNP Q05320
C	663	THR	-	expression tag	UNP Q05320
C	664	HIS	-	expression tag	UNP Q05320
C	665	HIS	-	expression tag	UNP Q05320
C	666	HIS	-	expression tag	UNP Q05320
C	667	HIS	-	expression tag	UNP Q05320
C	668	HIS	-	expression tag	UNP Q05320
C	669	HIS	-	expression tag	UNP Q05320
E	633	GLY	-	expression tag	UNP Q05320
E	634	SER	-	expression tag	UNP Q05320
E	635	GLY	-	expression tag	UNP Q05320
E	636	TYR	-	expression tag	UNP Q05320
E	637	ILE	-	expression tag	UNP Q05320
E	638	PRO	-	expression tag	UNP Q05320
E	639	GLU	-	expression tag	UNP Q05320
E	640	ALA	-	expression tag	UNP Q05320
E	641	PRO	-	expression tag	UNP Q05320
E	642	ARG	-	expression tag	UNP Q05320
E	643	ASP	-	expression tag	UNP Q05320
E	644	GLY	-	expression tag	UNP Q05320
E	645	GLN	-	expression tag	UNP Q05320
E	646	ALA	-	expression tag	UNP Q05320
E	647	TYR	-	expression tag	UNP Q05320
E	648	VAL	-	expression tag	UNP Q05320
E	649	ARG	-	expression tag	UNP Q05320
E	650	LYS	-	expression tag	UNP Q05320
E	651	ASP	-	expression tag	UNP Q05320
E	652	GLY	-	expression tag	UNP Q05320
E	653	GLU	-	expression tag	UNP Q05320
E	654	TRP	-	expression tag	UNP Q05320
E	655	VAL	-	expression tag	UNP Q05320
E	656	LEU	-	expression tag	UNP Q05320
E	657	LEU	-	expression tag	UNP Q05320
E	658	SER	-	expression tag	UNP Q05320
E	659	THR	-	expression tag	UNP Q05320
E	660	PHE	-	expression tag	UNP Q05320
E	661	LEU	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
E	662	GLY	-	expression tag	UNP Q05320
E	663	THR	-	expression tag	UNP Q05320
E	664	HIS	-	expression tag	UNP Q05320
E	665	HIS	-	expression tag	UNP Q05320
E	666	HIS	-	expression tag	UNP Q05320
E	667	HIS	-	expression tag	UNP Q05320
E	668	HIS	-	expression tag	UNP Q05320
E	669	HIS	-	expression tag	UNP Q05320
F	633	GLY	-	expression tag	UNP Q05320
F	634	SER	-	expression tag	UNP Q05320
F	635	GLY	-	expression tag	UNP Q05320
F	636	TYR	-	expression tag	UNP Q05320
F	637	ILE	-	expression tag	UNP Q05320
F	638	PRO	-	expression tag	UNP Q05320
F	639	GLU	-	expression tag	UNP Q05320
F	640	ALA	-	expression tag	UNP Q05320
F	641	PRO	-	expression tag	UNP Q05320
F	642	ARG	-	expression tag	UNP Q05320
F	643	ASP	-	expression tag	UNP Q05320
F	644	GLY	-	expression tag	UNP Q05320
F	645	GLN	-	expression tag	UNP Q05320
F	646	ALA	-	expression tag	UNP Q05320
F	647	TYR	-	expression tag	UNP Q05320
F	648	VAL	-	expression tag	UNP Q05320
F	649	ARG	-	expression tag	UNP Q05320
F	650	LYS	-	expression tag	UNP Q05320
F	651	ASP	-	expression tag	UNP Q05320
F	652	GLY	-	expression tag	UNP Q05320
F	653	GLU	-	expression tag	UNP Q05320
F	654	TRP	-	expression tag	UNP Q05320
F	655	VAL	-	expression tag	UNP Q05320
F	656	LEU	-	expression tag	UNP Q05320
F	657	LEU	-	expression tag	UNP Q05320
F	658	SER	-	expression tag	UNP Q05320
F	659	THR	-	expression tag	UNP Q05320
F	660	PHE	-	expression tag	UNP Q05320
F	661	LEU	-	expression tag	UNP Q05320
F	662	GLY	-	expression tag	UNP Q05320
F	663	THR	-	expression tag	UNP Q05320
F	664	HIS	-	expression tag	UNP Q05320
F	665	HIS	-	expression tag	UNP Q05320
F	666	HIS	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
F	667	HIS	-	expression tag	UNP Q05320
F	668	HIS	-	expression tag	UNP Q05320
F	669	HIS	-	expression tag	UNP Q05320

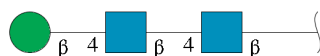
- Molecule 2 is a protein called Envelope glycoprotein, GP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	175	Total	C	N	O	S	0	0
			1277	818	225	229	5		
2	B	175	Total	C	N	O	S	0	0
			1271	812	223	231	5		
2	D	175	Total	C	N	O	S	0	0
			1273	813	223	233	4		

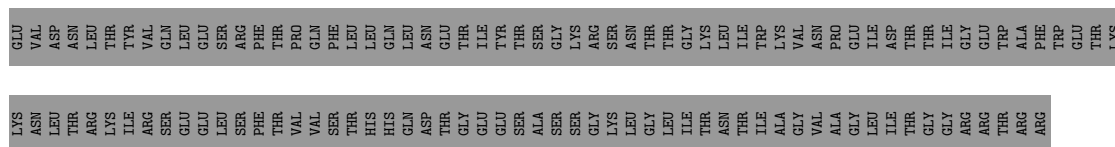
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	GLY	-	expression tag	UNP Q05320
A	31	SER	-	expression tag	UNP Q05320
A	42	ALA	THR	conflict	UNP Q05320
B	30	GLY	-	expression tag	UNP Q05320
B	31	SER	-	expression tag	UNP Q05320
B	42	ALA	THR	conflict	UNP Q05320
D	30	GLY	-	expression tag	UNP Q05320
D	31	SER	-	expression tag	UNP Q05320
D	42	ALA	THR	conflict	UNP Q05320

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

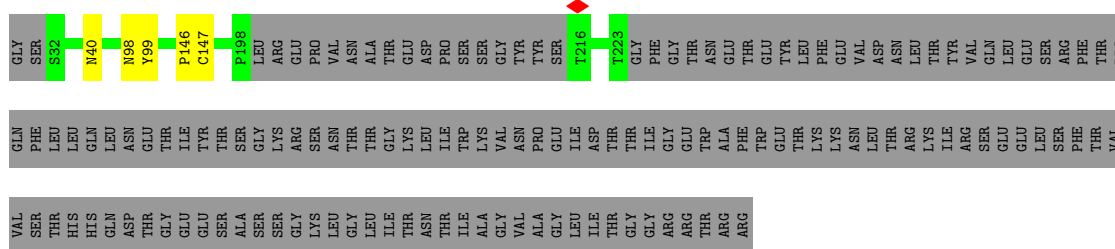


Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	3	Total	C	N	O		0	0
			39	22	2	15			
3	H	3	Total	C	N	O		0	0
			39	22	2	15			
3	I	3	Total	C	N	O		0	0
			39	22	2	15			



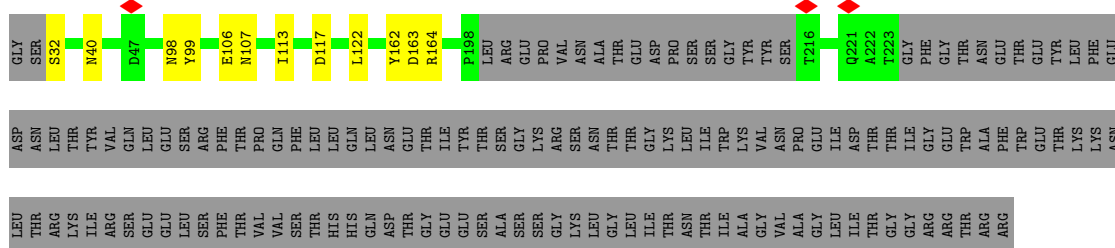
• Molecule 2: Envelope glycoprotein,GP1

Chain B: 53% 45%



• Molecule 2: Envelope glycoprotein,GP1

Chain D: 51% 45%



• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 33% 67%



• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 33% 67%



• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I: 100%

MAG1
MAG2
BMA3

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	189700	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 ($e^-/\text{\AA}^2$)	56.69	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.145	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.032	Depositor
Recommended contour level	0.0213	Depositor
Map size (Å)	215.04, 215.04, 215.04	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.84, 0.84, 0.84	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, BMA

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	C	0.19	0/767	0.37	0/1045
1	E	0.19	0/767	0.31	0/1045
1	F	0.19	0/770	0.32	0/1049
2	A	0.18	0/1308	0.33	0/1781
2	B	0.19	0/1302	0.31	0/1775
2	D	0.19	0/1304	0.31	0/1778
All	All	0.19	0/6218	0.32	0/8473

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	C	749	0	741	3	0
1	E	749	0	739	1	0
1	F	752	0	743	2	0
2	A	1277	0	1214	8	0
2	B	1271	0	1189	3	0
2	D	1273	0	1188	9	0
3	G	39	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	39	0	34	0	0
3	I	39	0	34	0	0
All	All	6188	0	5916	23	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 2.

All (23) 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)
2:D:106:GLU:OE1	2:D:107:ASN:ND2	2.31	0.64
2:A:56:LYS:NZ	2:A:188:GLN:OE1	2.31	0.62
2:D:113:ILE:HG22	2:D:122:LEU:HD11	1.84	0.59
2:B:98:ASN:OD1	2:B:99:TYR:N	2.37	0.58
2:A:117:ASP:C	2:A:117:ASP:OD1	2.51	0.54
2:A:61:ASN:OD1	2:A:61:ASN:C	2.51	0.52
2:D:117:ASP:C	2:D:117:ASP:OD1	2.53	0.52
2:B:40:ASN:O	2:B:40:ASN:ND2	2.42	0.51
2:B:146:PRO:O	2:B:147:CYS:C	2.55	0.49
1:E:552:ASP:OD1	1:E:553:GLY:N	2.46	0.48
1:F:523:GLU:OE1	2:A:32:SER:OG	2.25	0.47
1:C:596:ARG:O	1:C:597:TRP:HE3	1.98	0.47
2:D:32:SER:O	2:D:32:SER:OG	2.28	0.45
2:D:98:ASN:OD1	2:D:99:TYR:N	2.48	0.45
1:C:562:ALA:HB1	2:A:180:VAL:HG23	1.99	0.44
1:F:517:TYR:H	1:F:517:TYR:HD1	1.67	0.42
2:D:163:ASP:OD1	2:D:164:ARG:N	2.46	0.42
2:D:162:TYR:O	2:D:163:ASP:C	2.62	0.42
2:D:40:ASN:O	2:D:40:ASN:ND2	2.53	0.41
2:D:113:ILE:HD12	2:D:113:ILE:HA	1.91	0.41
2:A:111:LEU:HD23	2:A:141:VAL:HB	2.03	0.41
1:C:516:HIS:O	2:A:103:GLU:HA	2.21	0.40
2:A:115:LYS:HG3	2:A:121:CYS:SG	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	C	93/168 (55%)	89 (96%)	4 (4%)	0	100	100
1	E	93/168 (55%)	92 (99%)	1 (1%)	0	100	100
1	F	93/168 (55%)	90 (97%)	3 (3%)	0	100	100
2	A	171/321 (53%)	164 (96%)	7 (4%)	0	100	100
2	B	171/321 (53%)	167 (98%)	4 (2%)	0	100	100
2	D	171/321 (53%)	166 (97%)	5 (3%)	0	100	100
All	All	792/1467 (54%)	768 (97%)	24 (3%)	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	C	75/140 (54%)	74 (99%)	1 (1%)	61	74
1	E	75/140 (54%)	75 (100%)	0	100	100
1	F	76/140 (54%)	76 (100%)	0	100	100
2	A	125/271 (46%)	124 (99%)	1 (1%)	73	79
2	B	123/271 (45%)	123 (100%)	0	100	100
2	D	123/271 (45%)	123 (100%)	0	100	100
All	All	597/1233 (48%)	595 (100%)	2 (0%)	84	85

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	597	TRP
2	A	77	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	508	GLN
2	A	40	ASN
2	A	98	ASN
2	D	73	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

9 monosaccharides 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	$\# Z > 2$
3	NAG	G	1	1,3	14,14,15	0.80	0	17,19,21	0.91	1 (5%)
3	NAG	G	2	3	14,14,15	0.76	0	17,19,21	0.87	0
3	BMA	G	3	3	11,11,12	0.87	0	15,15,17	2.75	7 (46%)
3	NAG	H	1	1,3	14,14,15	0.81	0	17,19,21	0.91	1 (5%)
3	NAG	H	2	3	14,14,15	0.73	0	17,19,21	0.88	0
3	BMA	H	3	3	11,11,12	0.85	0	15,15,17	2.83	6 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	I	1	1,3	14,14,15	0.78	0	17,19,21	0.94	1 (5%)
3	NAG	I	2	3	14,14,15	0.76	0	17,19,21	0.95	1 (5%)
3	BMA	I	3	3	11,11,12	0.86	0	15,15,17	2.71	6 (40%)

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
3	NAG	G	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	BMA	G	3	3	-	0/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
3	NAG	I	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	BMA	I	3	3	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	3	BMA	C1-O5-C5	8.02	122.93	112.19
3	I	3	BMA	C1-O5-C5	7.53	122.28	112.19
3	G	3	BMA	C1-O5-C5	7.50	122.24	112.19
3	H	3	BMA	C3-C4-C5	4.17	117.79	110.23
3	G	3	BMA	C3-C4-C5	4.15	117.75	110.23
3	I	3	BMA	C3-C4-C5	4.06	117.59	110.23
3	G	3	BMA	C2-C3-C4	3.75	117.46	110.86
3	I	3	BMA	C2-C3-C4	3.57	117.15	110.86
3	H	3	BMA	C2-C3-C4	3.51	117.04	110.86
3	G	3	BMA	O3-C3-C2	-2.54	104.87	110.05
3	G	1	NAG	O5-C1-C2	-2.52	107.39	111.29
3	I	3	BMA	O3-C3-C2	-2.51	104.94	110.05
3	H	3	BMA	O4-C4-C3	-2.48	104.52	110.38
3	G	3	BMA	O4-C4-C3	-2.47	104.56	110.38
3	I	3	BMA	O5-C5-C4	2.47	116.83	110.83
3	H	3	BMA	O3-C3-C2	-2.46	105.03	110.05
3	H	3	BMA	O5-C5-C4	2.46	116.80	110.83

Continued on next page...

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	2	NAG	C1-O5-C5	2.44	115.46	112.19
3	I	1	NAG	O5-C1-C2	-2.42	107.55	111.29
3	G	3	BMA	O5-C5-C4	2.30	116.43	110.83
3	H	1	NAG	O5-C1-C2	-2.29	107.74	111.29
3	I	3	BMA	O4-C4-C3	-2.25	105.06	110.38
3	G	3	BMA	C1-C2-C3	2.14	112.75	109.64

There are no chirality outliers.

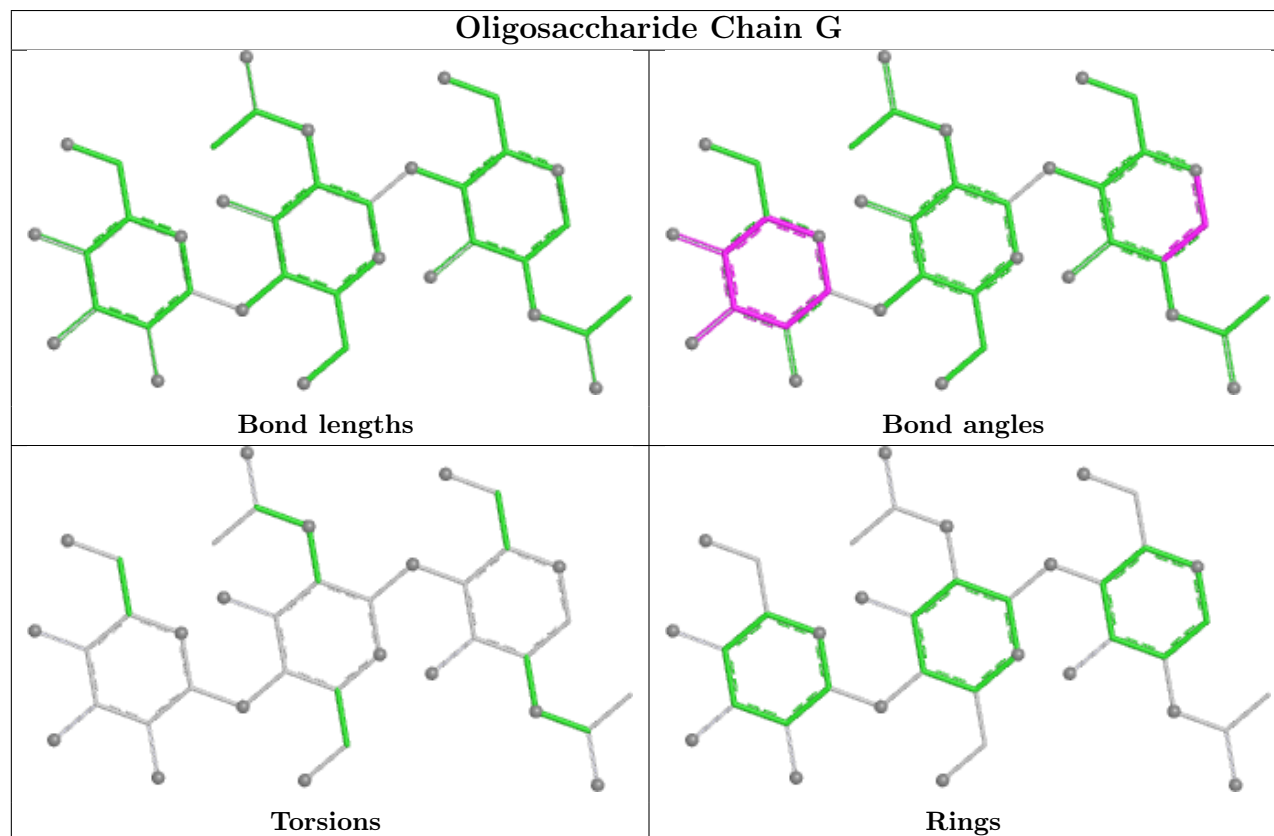
All (1) torsion outliers are listed below:

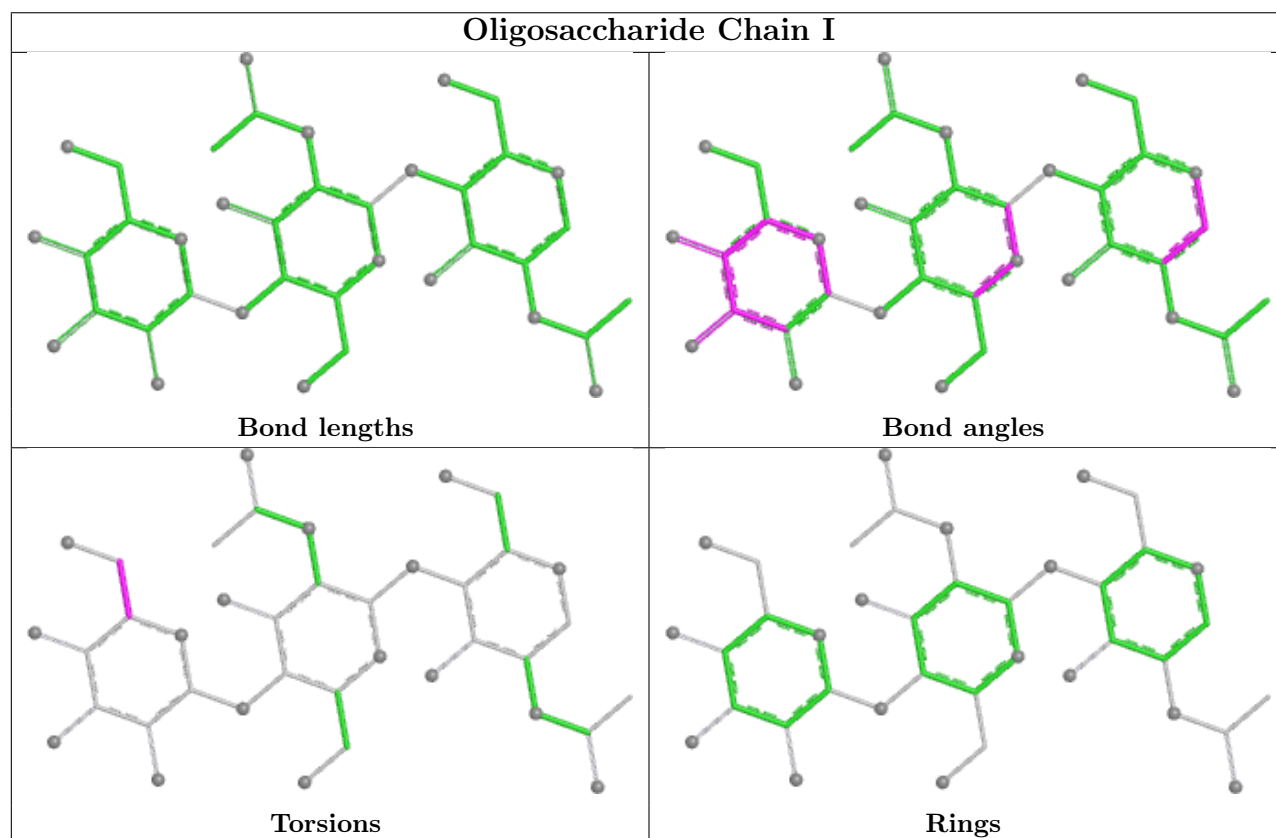
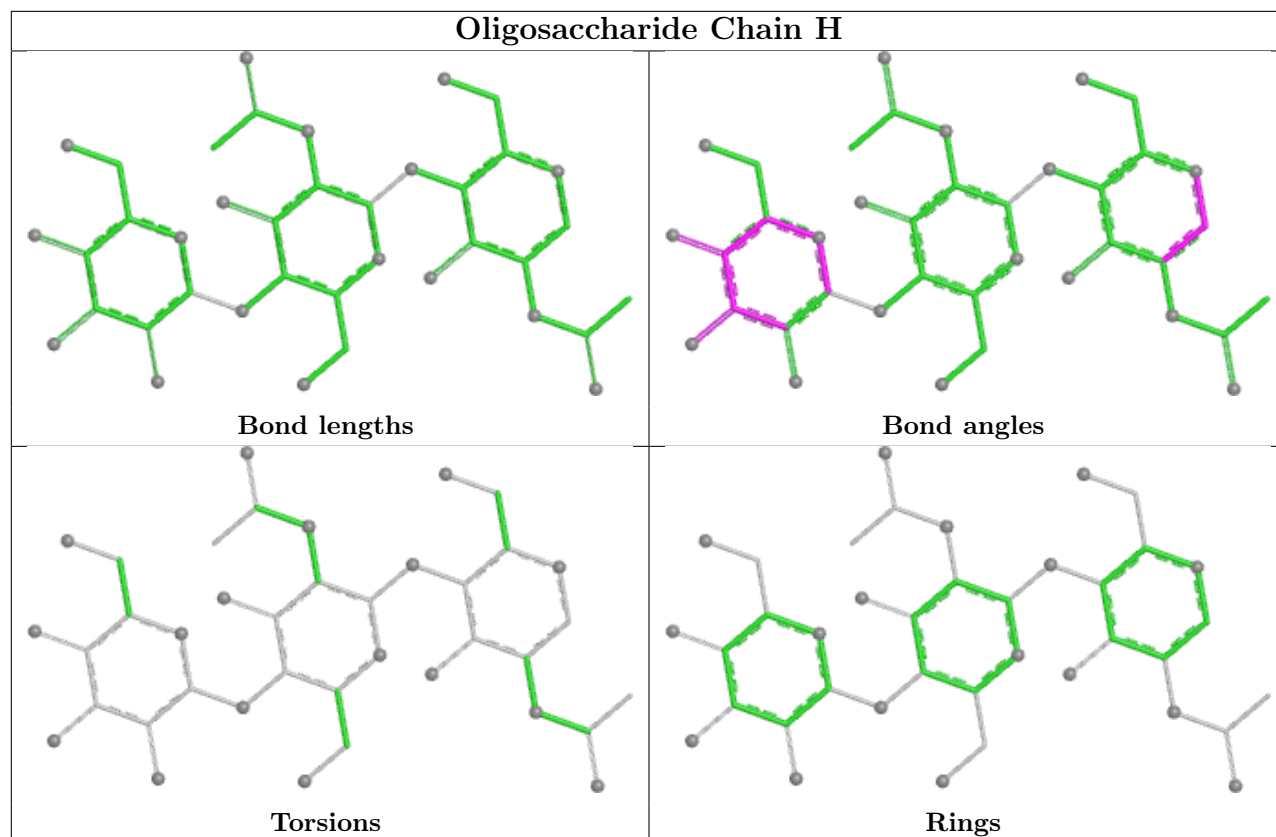
Mol	Chain	Res	Type	Atoms
3	I	3	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

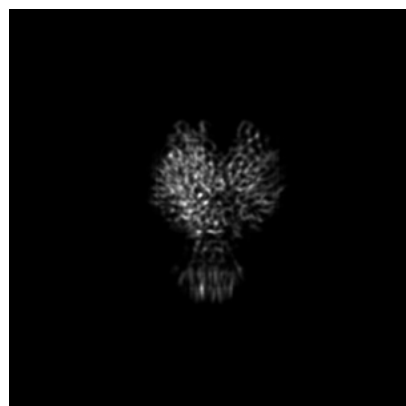
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-48271. These allow visual inspection of the internal detail of the map and identification of artifacts.

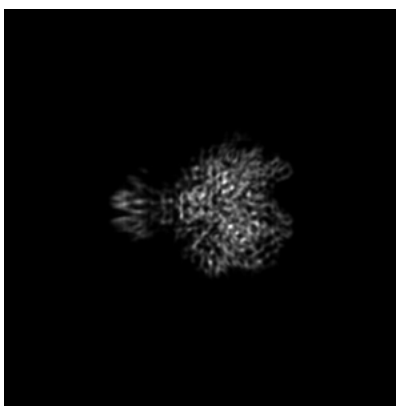
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

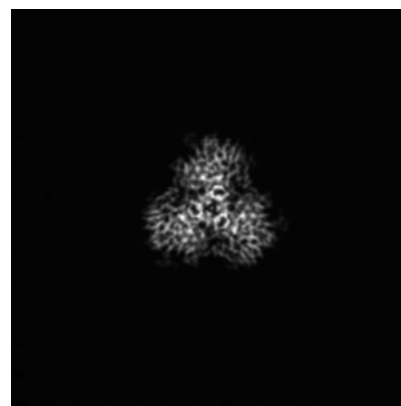
6.1.1 Primary map



X

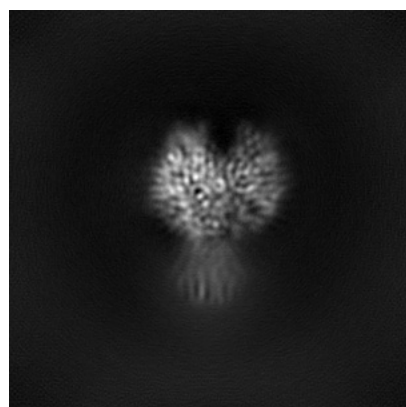


Y

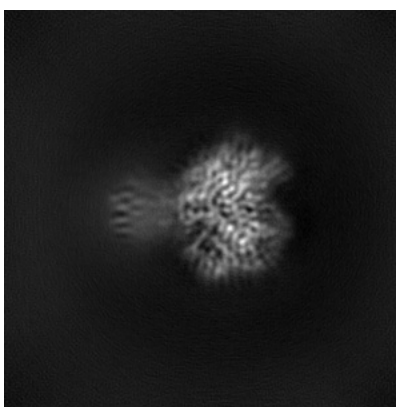


Z

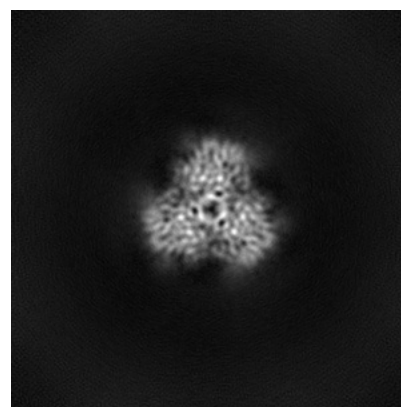
6.1.2 Raw map



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

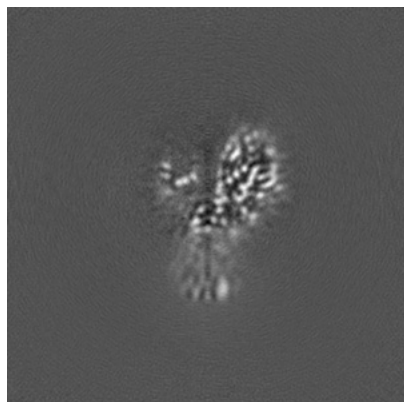


Y Index: 128

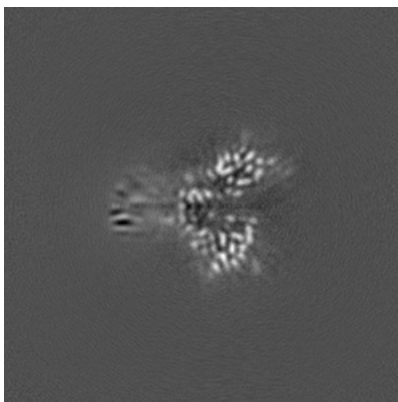


Z Index: 128

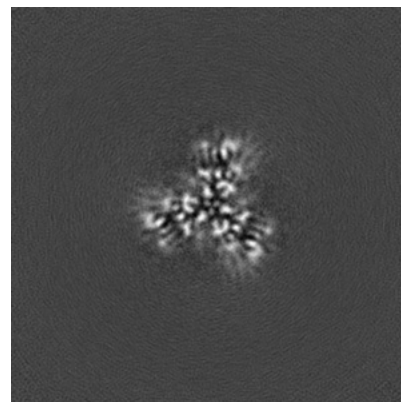
6.2.2 Raw map



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: 131

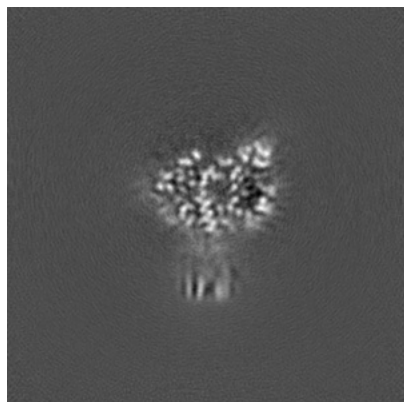


Y Index: 123

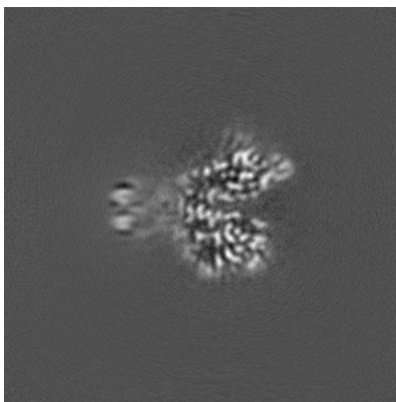


Z Index: 140

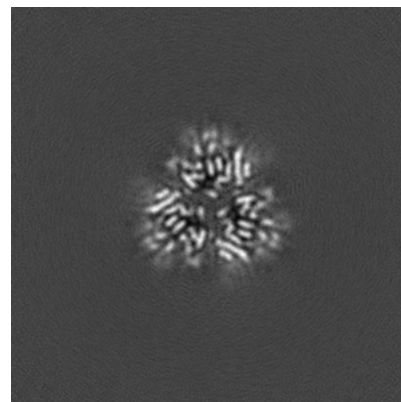
6.3.2 Raw map



X Index: 136



Y Index: 123

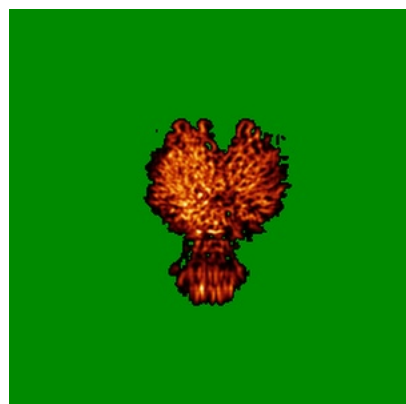


Z Index: 141

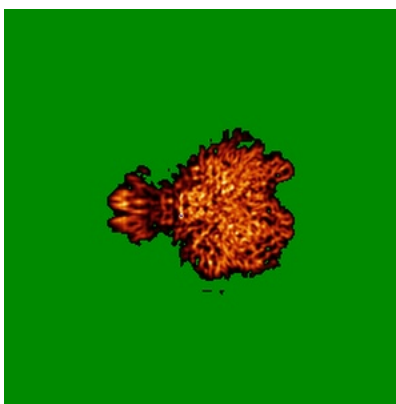
The images above show the largest variance slices of the map in three orthogonal directions.

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

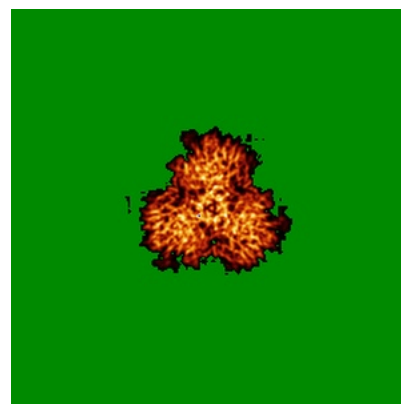
6.4.1 Primary map



X

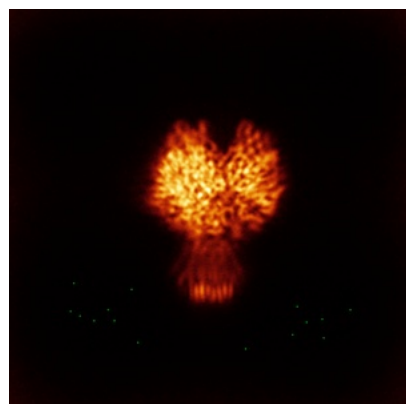


Y

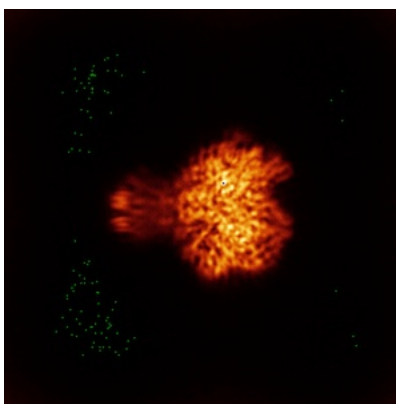


Z

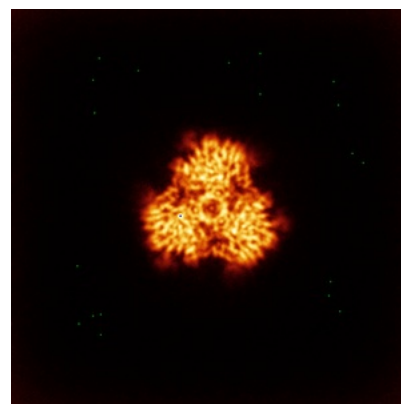
6.4.2 Raw map



X



Y

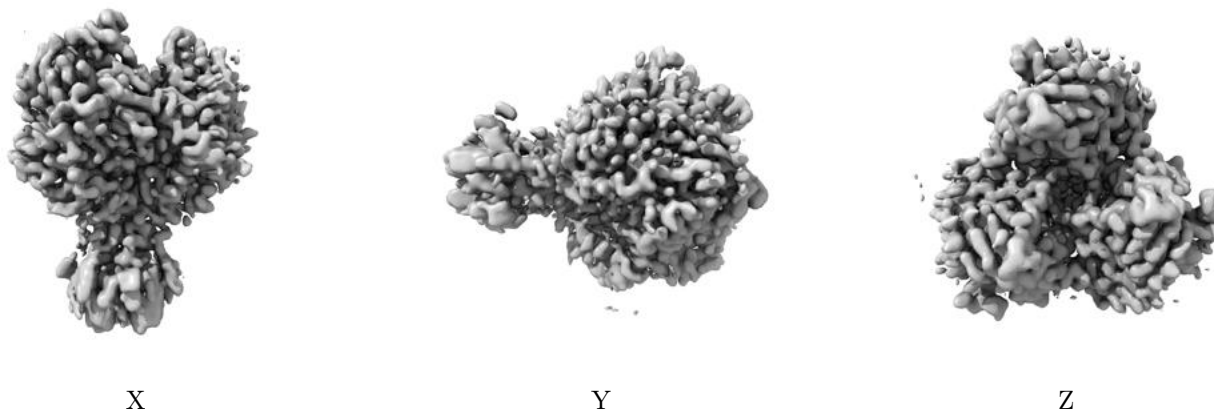


Z

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.

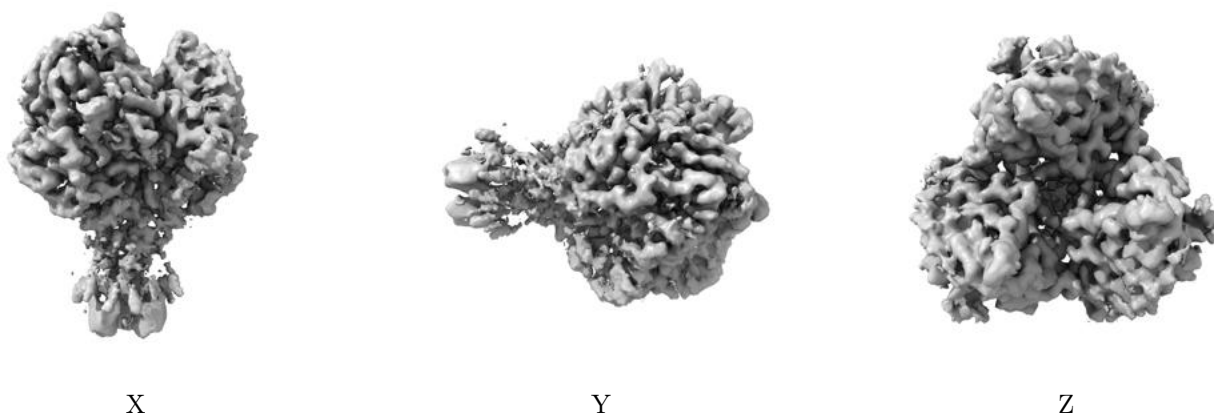
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0213. 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.

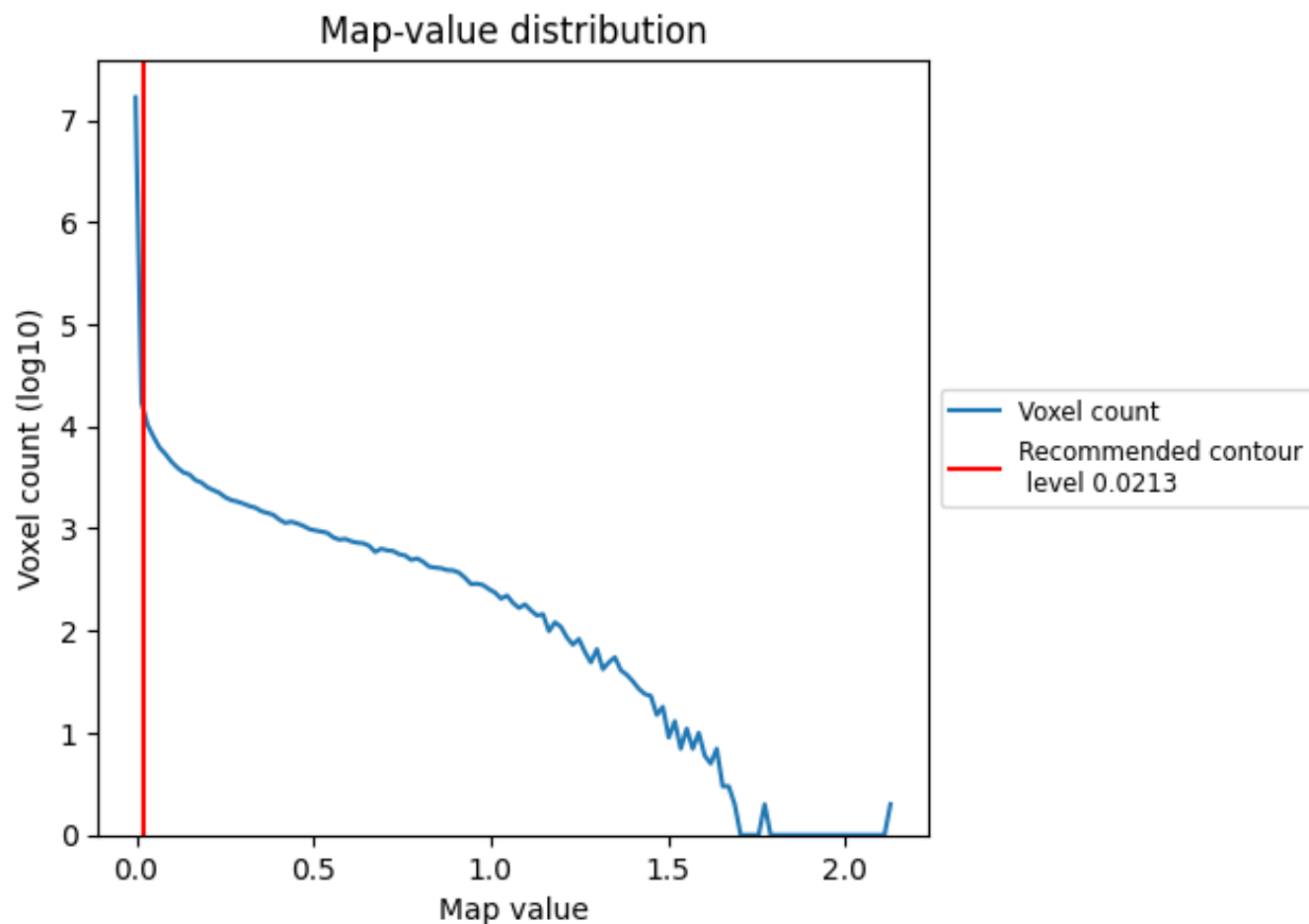
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

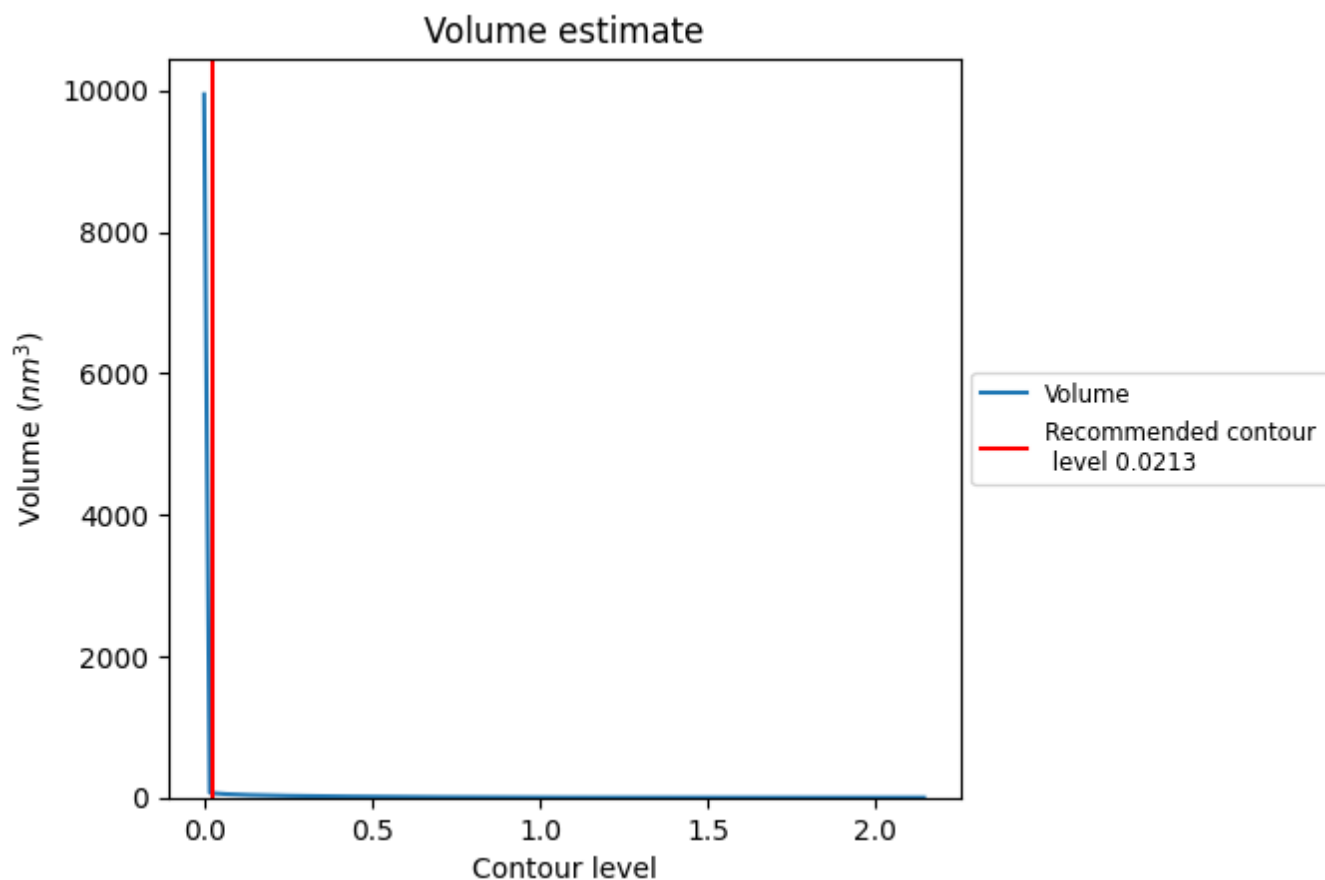
This section contains the results of statistical analysis of the map.

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.

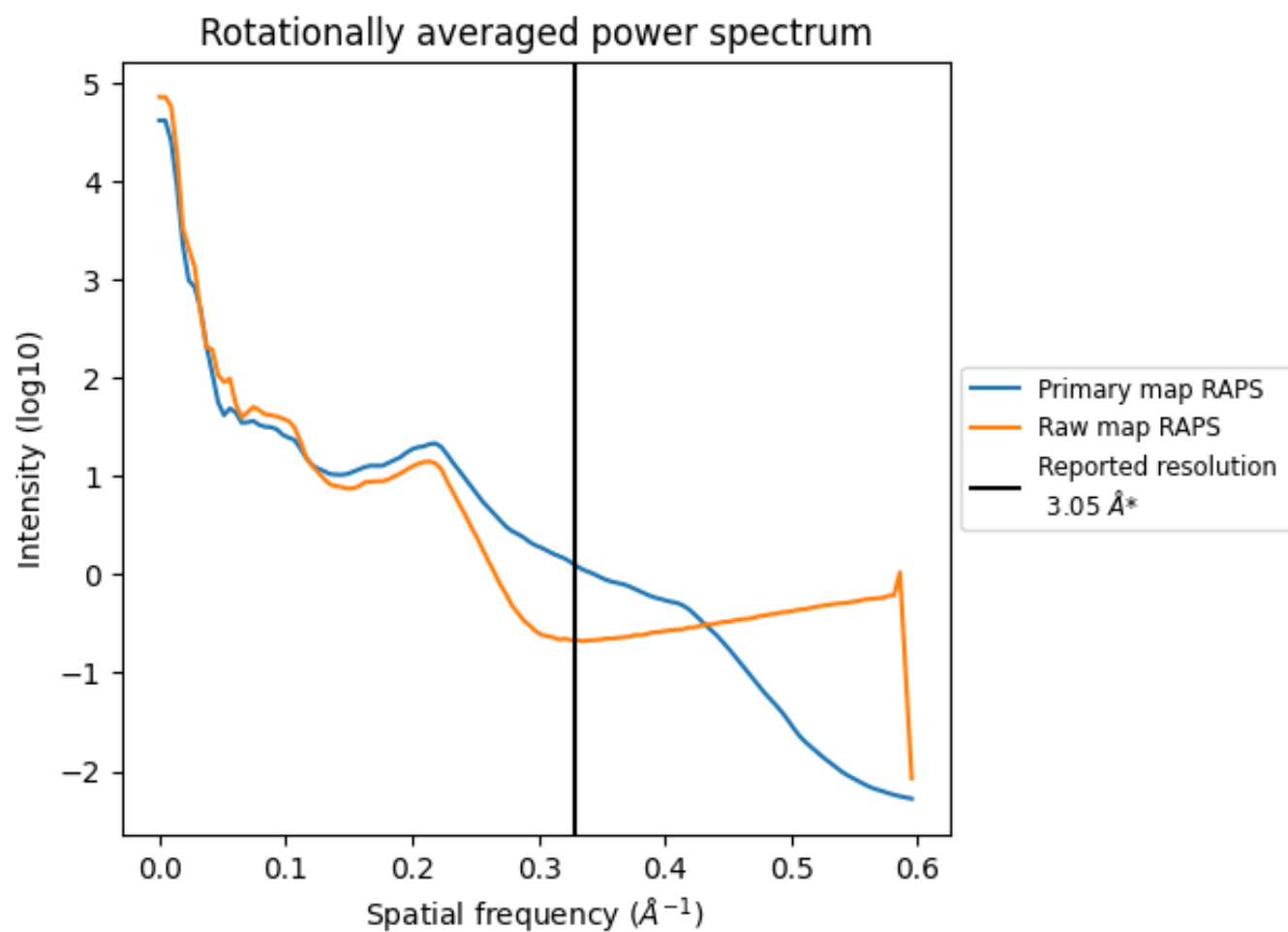
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 65 nm^3 ; this corresponds to an approximate mass of 59 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 ⓘ

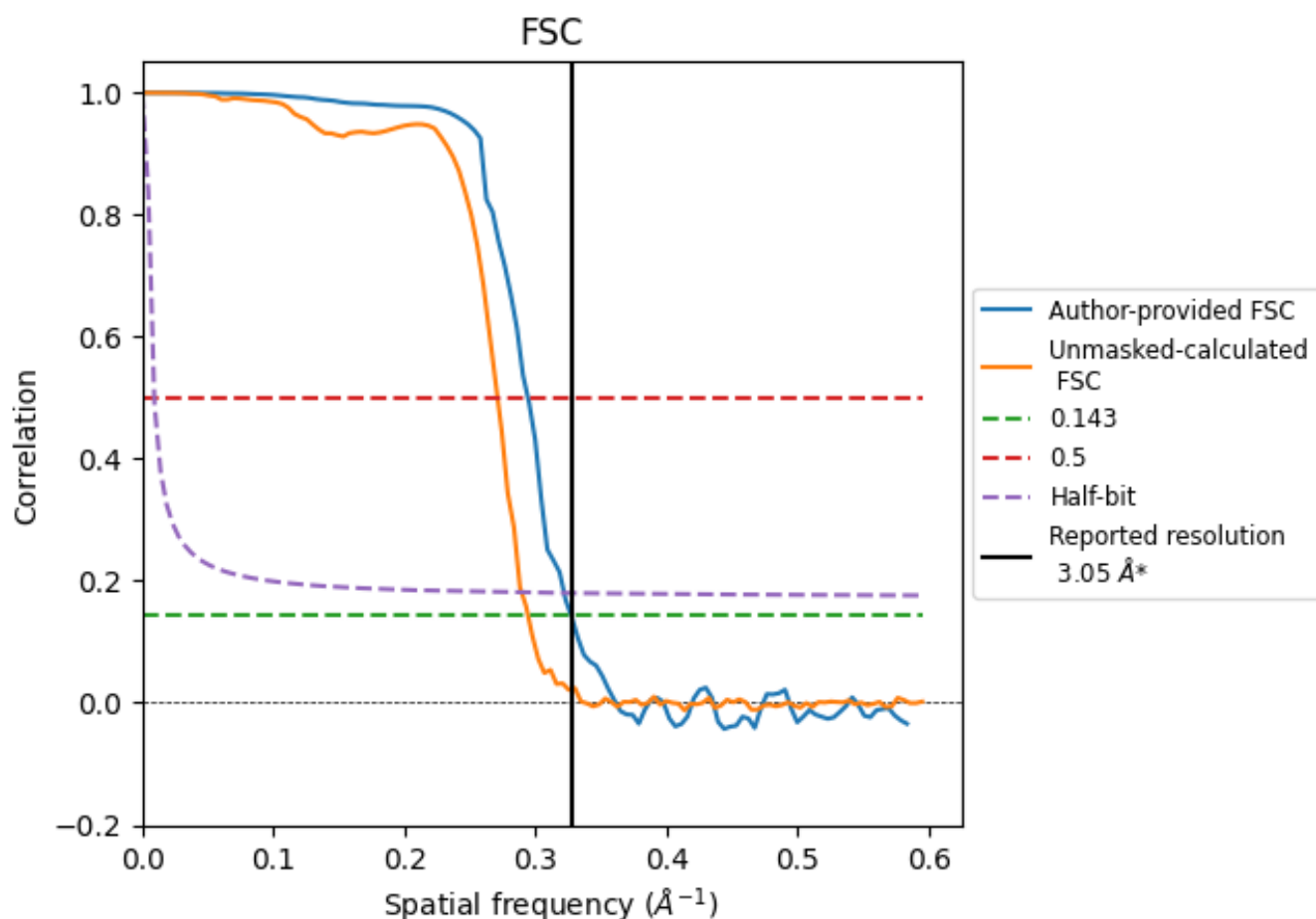


*Reported resolution corresponds to spatial frequency of 0.328 Å⁻¹

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 0.328 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

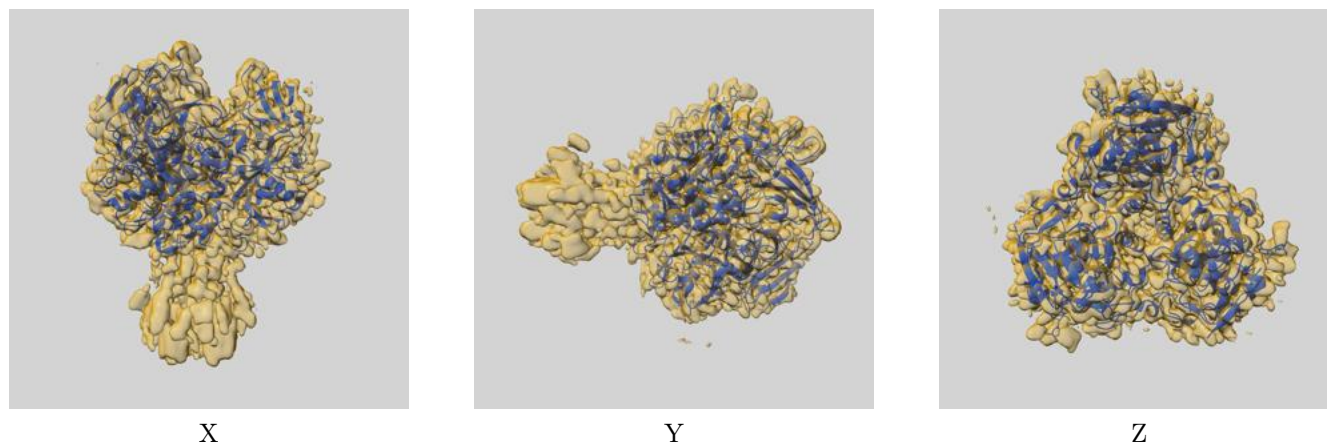
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.05	-	-
Author-provided FSC curve	3.05	3.40	3.11
Unmasked-calculated*	3.40	3.69	3.45

*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 3.40 differs from the reported value 3.05 by more than 10 %

9 Map-model fit [i](#)

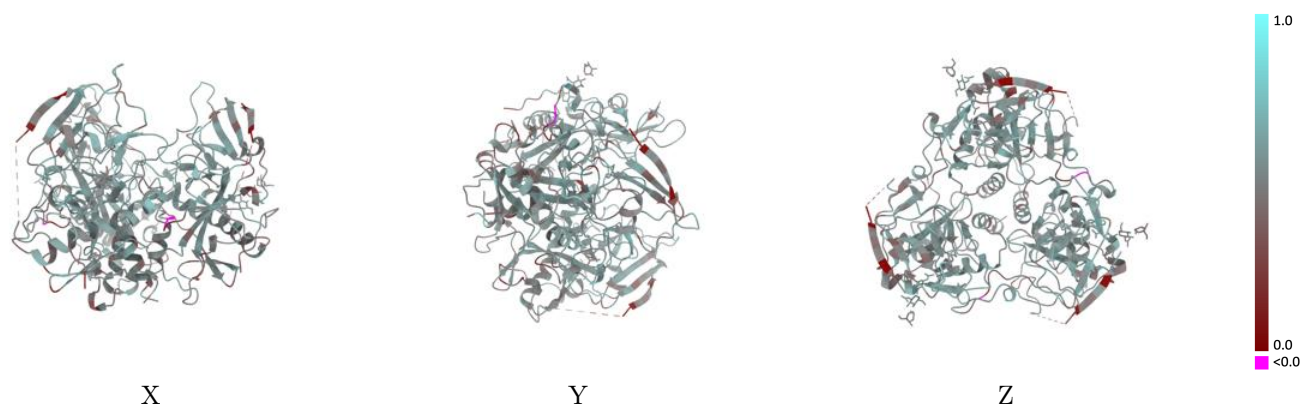
This section contains information regarding the fit between EMDB map EMD-48271 and PDB model 9MHA. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



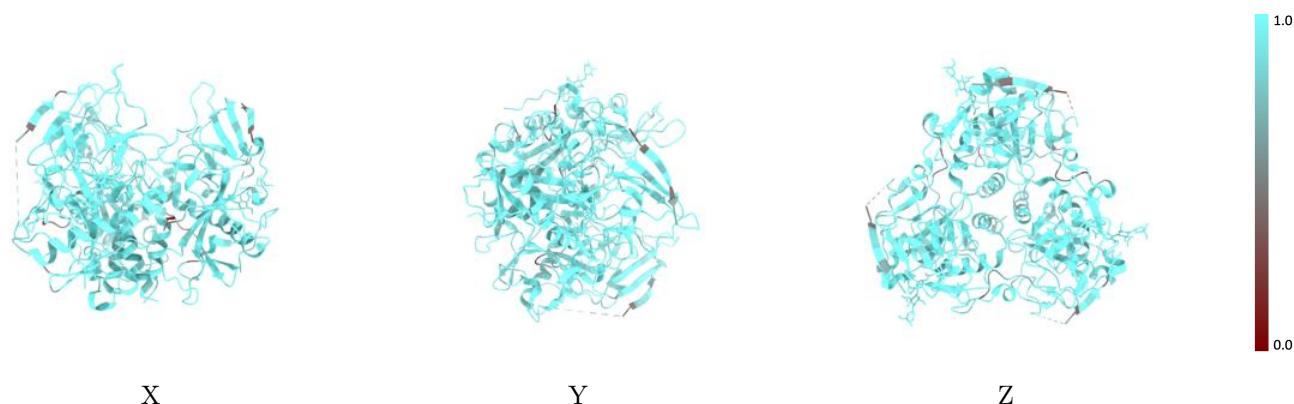
The images above show the 3D surface view of the map at the recommended contour level 0.0213 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



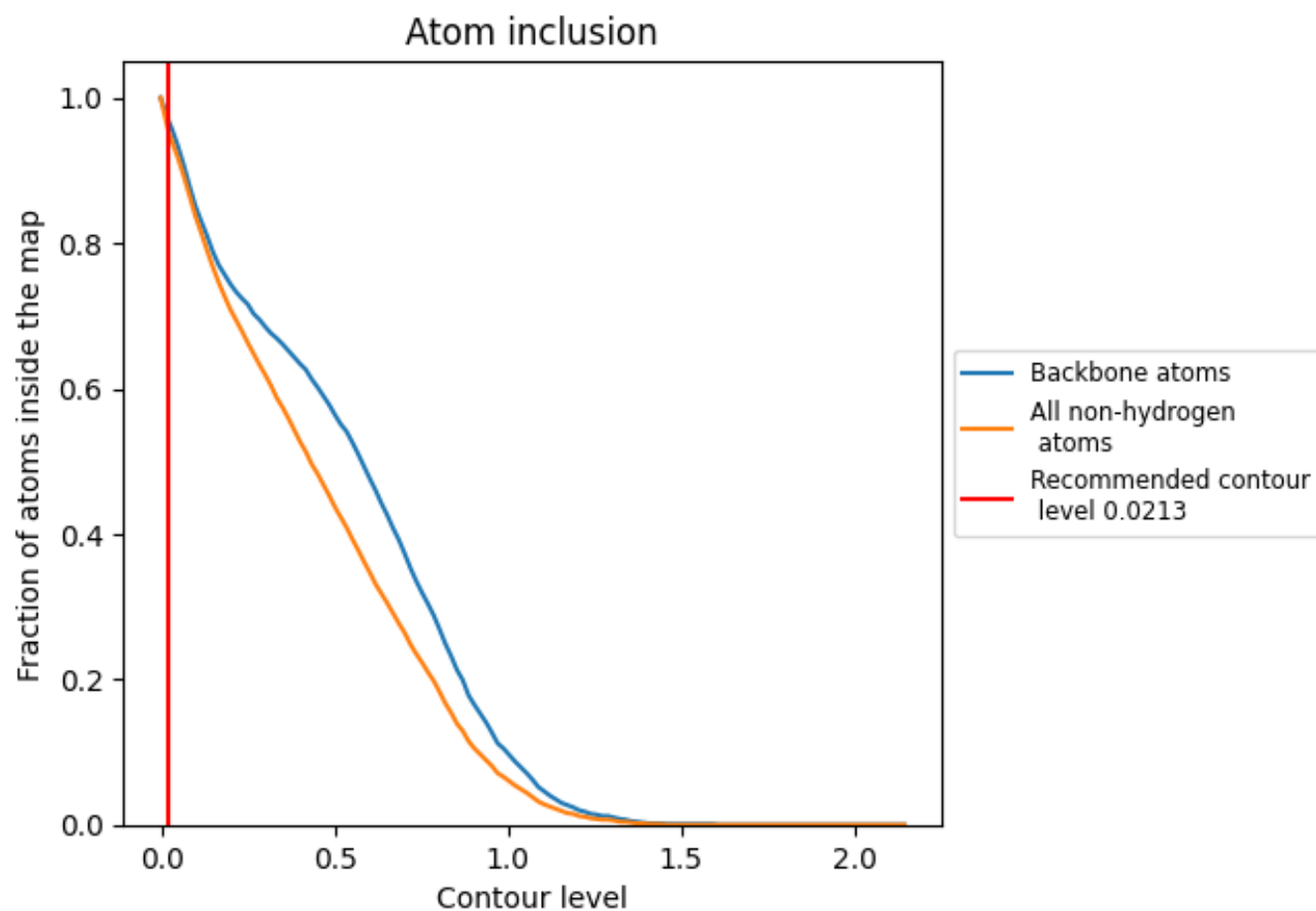
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0213).

9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 95% 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 (0.0213) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9530	0.5140
A	0.9690	0.5270
B	0.9620	0.5220
C	0.9360	0.5000
D	0.9500	0.5150
E	0.9450	0.5060
F	0.9400	0.4990
G	0.9490	0.5000
H	0.9490	0.5320
I	0.9740	0.5390

1.0

0.0

<0.0