



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 03:11 PM UTC

PDB ID : 9BSV / pdb_00009bsv
EMDB ID : EMD-44873
Title : EBOV GP/Nanosota-EB2 complex
Authors : Bu, F.; Ye, G.; Liu, B.; Li, F.
Deposited on : 2024-05-13
Resolution : 3.10 Å(reported)
Based on initial model : 5jq7

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:

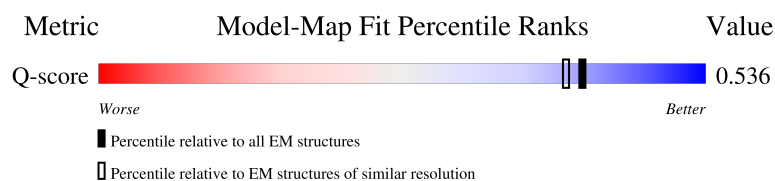
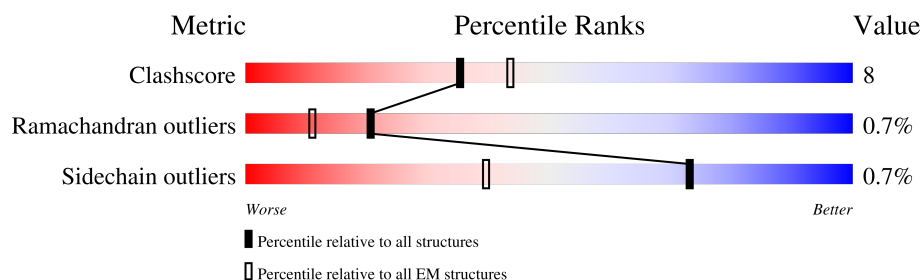
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
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 3.10 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.



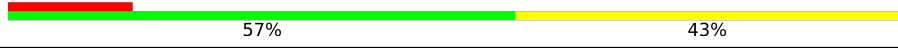
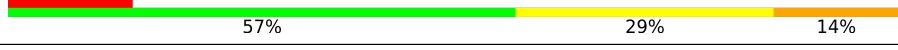
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14724 (2.60 - 3.60)

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	D	144	
1	E	144	
1	F	144	
2	A	706	

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Mol	Chain	Length	Quality of chain
2	B	706	
2	C	706	
3	G	7	
4	H	7	
5	I	6	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	MAN	I	4	-	-	X	-
6	MAN	E	201	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9879 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 Nanosota-EB2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	D	123	Total	C	N	O	S	0	0
			938	580	168	186	4		
1	E	123	Total	C	N	O	S	0	0
			938	580	168	186	4		
1	F	123	Total	C	N	O	S	0	0
			938	580	168	186	4		

- Molecule 2 is a protein called Envelope glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	293	Total	C	N	O	S	0	0
			2272	1450	395	416	11		
2	B	293	Total	C	N	O	S	0	0
			2272	1450	395	416	11		
2	C	293	Total	C	N	O	S	0	0
			2272	1450	395	416	11		

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	677	GLY	-	expression tag	UNP Q05320
A	678	SER	-	expression tag	UNP Q05320
A	679	GLY	-	expression tag	UNP Q05320
A	680	TYR	-	expression tag	UNP Q05320
A	681	ILE	-	expression tag	UNP Q05320
A	682	PRO	-	expression tag	UNP Q05320
A	683	GLU	-	expression tag	UNP Q05320
A	684	ALA	-	expression tag	UNP Q05320
A	685	PRO	-	expression tag	UNP Q05320
A	686	ARG	-	expression tag	UNP Q05320
A	687	ASP	-	expression tag	UNP Q05320
A	688	GLY	-	expression tag	UNP Q05320
A	689	GLN	-	expression tag	UNP Q05320

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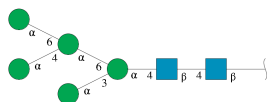
Chain	Residue	Modelled	Actual	Comment	Reference
A	690	ALA	-	expression tag	UNP Q05320
A	691	TYR	-	expression tag	UNP Q05320
A	692	VAL	-	expression tag	UNP Q05320
A	693	ARG	-	expression tag	UNP Q05320
A	694	LYS	-	expression tag	UNP Q05320
A	695	ASP	-	expression tag	UNP Q05320
A	696	GLY	-	expression tag	UNP Q05320
A	697	GLU	-	expression tag	UNP Q05320
A	698	TRP	-	expression tag	UNP Q05320
A	699	VAL	-	expression tag	UNP Q05320
A	700	LEU	-	expression tag	UNP Q05320
A	701	LEU	-	expression tag	UNP Q05320
A	702	SER	-	expression tag	UNP Q05320
A	703	THR	-	expression tag	UNP Q05320
A	704	PHE	-	expression tag	UNP Q05320
A	705	LEU	-	expression tag	UNP Q05320
A	706	GLY	-	expression tag	UNP Q05320
B	677	GLY	-	expression tag	UNP Q05320
B	678	SER	-	expression tag	UNP Q05320
B	679	GLY	-	expression tag	UNP Q05320
B	680	TYR	-	expression tag	UNP Q05320
B	681	ILE	-	expression tag	UNP Q05320
B	682	PRO	-	expression tag	UNP Q05320
B	683	GLU	-	expression tag	UNP Q05320
B	684	ALA	-	expression tag	UNP Q05320
B	685	PRO	-	expression tag	UNP Q05320
B	686	ARG	-	expression tag	UNP Q05320
B	687	ASP	-	expression tag	UNP Q05320
B	688	GLY	-	expression tag	UNP Q05320
B	689	GLN	-	expression tag	UNP Q05320
B	690	ALA	-	expression tag	UNP Q05320
B	691	TYR	-	expression tag	UNP Q05320
B	692	VAL	-	expression tag	UNP Q05320
B	693	ARG	-	expression tag	UNP Q05320
B	694	LYS	-	expression tag	UNP Q05320
B	695	ASP	-	expression tag	UNP Q05320
B	696	GLY	-	expression tag	UNP Q05320
B	697	GLU	-	expression tag	UNP Q05320
B	698	TRP	-	expression tag	UNP Q05320
B	699	VAL	-	expression tag	UNP Q05320
B	700	LEU	-	expression tag	UNP Q05320
B	701	LEU	-	expression tag	UNP Q05320

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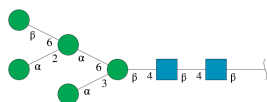
Chain	Residue	Modelled	Actual	Comment	Reference
B	702	SER	-	expression tag	UNP Q05320
B	703	THR	-	expression tag	UNP Q05320
B	704	PHE	-	expression tag	UNP Q05320
B	705	LEU	-	expression tag	UNP Q05320
B	706	GLY	-	expression tag	UNP Q05320
C	677	GLY	-	expression tag	UNP Q05320
C	678	SER	-	expression tag	UNP Q05320
C	679	GLY	-	expression tag	UNP Q05320
C	680	TYR	-	expression tag	UNP Q05320
C	681	ILE	-	expression tag	UNP Q05320
C	682	PRO	-	expression tag	UNP Q05320
C	683	GLU	-	expression tag	UNP Q05320
C	684	ALA	-	expression tag	UNP Q05320
C	685	PRO	-	expression tag	UNP Q05320
C	686	ARG	-	expression tag	UNP Q05320
C	687	ASP	-	expression tag	UNP Q05320
C	688	GLY	-	expression tag	UNP Q05320
C	689	GLN	-	expression tag	UNP Q05320
C	690	ALA	-	expression tag	UNP Q05320
C	691	TYR	-	expression tag	UNP Q05320
C	692	VAL	-	expression tag	UNP Q05320
C	693	ARG	-	expression tag	UNP Q05320
C	694	LYS	-	expression tag	UNP Q05320
C	695	ASP	-	expression tag	UNP Q05320
C	696	GLY	-	expression tag	UNP Q05320
C	697	GLU	-	expression tag	UNP Q05320
C	698	TRP	-	expression tag	UNP Q05320
C	699	VAL	-	expression tag	UNP Q05320
C	700	LEU	-	expression tag	UNP Q05320
C	701	LEU	-	expression tag	UNP Q05320
C	702	SER	-	expression tag	UNP Q05320
C	703	THR	-	expression tag	UNP Q05320
C	704	PHE	-	expression tag	UNP Q05320
C	705	LEU	-	expression tag	UNP Q05320
C	706	GLY	-	expression tag	UNP Q05320

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



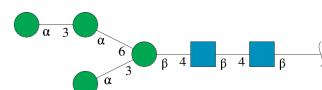
Mol	Chain	Residues	Atoms				AltConf	Trace
3	G	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-[beta-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



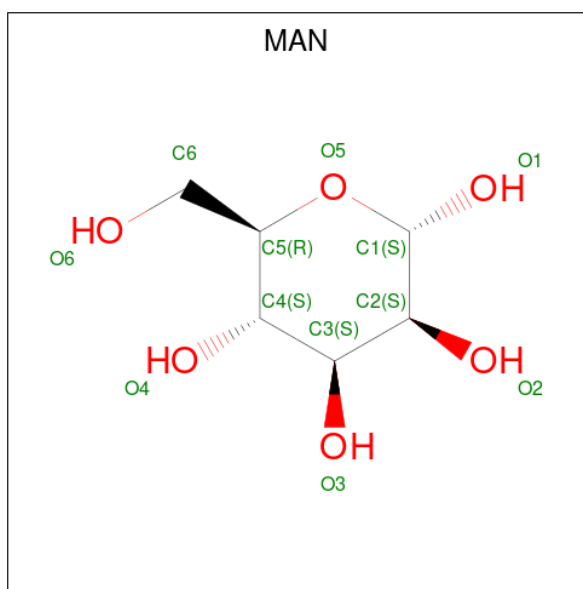
Mol	Chain	Residues	Atoms				AltConf	Trace
4	H	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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
5	I	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 6 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).



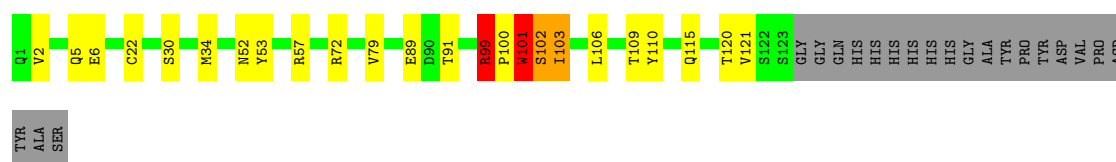
Mol	Chain	Residues	Atoms			AltConf
6	E	1	Total	C	O	0
			11	6	5	

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: Nanosota-EB2

Chain D: 



- Molecule 1: Nanosota-EB2

Chain E: 



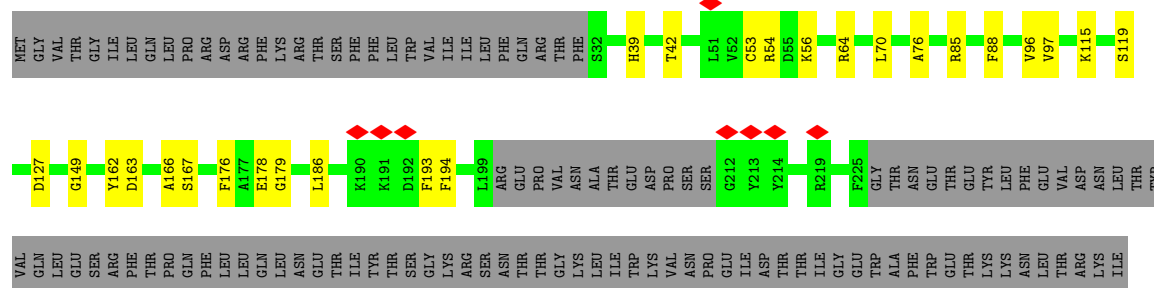
- Molecule 1: Nanosota-EB2

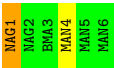
Chain F: 



- Molecule 2: Envelope glycoprotein

Chain A: 





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	404158	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.668	Depositor
Minimum map value	-0.360	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.079	Depositor
Map size ($\text{\AA}$)	283.30664, 283.30664, 283.30664	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.88533324, 0.88533324, 0.88533324	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, BMA, 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	D	0.31	0/957	0.53	1/1296 (0.1%)
1	E	0.26	0/957	0.41	0/1296
1	F	0.29	0/957	0.42	1/1296 (0.1%)
2	A	0.17	0/2330	0.35	0/3166
2	B	0.18	0/2330	0.34	0/3166
2	C	0.18	0/2330	0.36	0/3166
All	All	0.21	0/9861	0.38	2/13386 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	101	TRP	CA-C-O	-5.97	115.27	120.88
1	F	100	PRO	CA-C-O	-5.21	115.48	121.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	99	ARG	Sidechain

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	D	938	0	890	18	0
1	E	938	0	890	16	0
1	F	938	0	890	13	0
2	A	2272	0	2215	37	0
2	B	2272	0	2213	26	0
2	C	2272	0	2213	36	0
3	G	83	0	70	1	0
4	H	83	0	70	6	0
5	I	72	0	61	13	0
6	E	11	0	10	12	0
All	All	9879	0	9522	154	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 (154) 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)
6:E:201:MAN:C4	5:I:4:MAN:O6	1.72	1.36
2:A:521:GLN:OE1	2:A:544:ILE:HD11	1.28	1.27
2:A:521:GLN:OE1	2:A:544:ILE:CD1	1.82	1.26
6:E:201:MAN:C4	5:I:4:MAN:C6	2.19	1.20
6:E:201:MAN:C4	5:I:4:MAN:H61	1.89	0.99
6:E:201:MAN:C5	5:I:4:MAN:O6	2.13	0.96
6:E:201:MAN:C6	5:I:4:MAN:O6	2.13	0.96
6:E:201:MAN:O6	5:I:4:MAN:O6	1.83	0.96
6:E:201:MAN:C3	5:I:4:MAN:H61	1.96	0.94
6:E:201:MAN:H3	5:I:4:MAN:H61	1.56	0.88
2:C:52:VAL:H	2:C:596:ARG:HH21	1.29	0.81
2:A:521:GLN:NE2	2:A:522:ASP:OD1	2.20	0.73
2:A:85:ARG:NH1	2:A:178:GLU:OE2	2.20	0.73
2:A:521:GLN:OE1	2:A:544:ILE:HD12	1.88	0.73
6:E:201:MAN:C6	5:I:4:MAN:HO6	2.00	0.71
1:D:5:GLN:OE1	1:D:115:GLN:NE2	2.24	0.70
1:E:29:PHE:HE1	1:E:34:MET:HE2	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:5:GLN:OE1	1:F:115:GLN:NE2	2.25	0.69
2:C:551:GLN:OE1	2:C:551:GLN:N	2.26	0.68
1:E:52:ASN:O	1:E:72:ARG:NH1	2.27	0.68
2:B:52:VAL:HB	2:B:55:ASP:HB2	1.77	0.67
1:D:52:ASN:O	1:D:72:ARG:NH1	2.27	0.67
1:F:83:MET:HE2	1:F:86:LEU:HD21	1.77	0.66
1:D:89:GLU:OE1	1:D:89:GLU:N	2.29	0.66
1:E:5:GLN:OE1	1:E:115:GLN:NE2	2.28	0.66
2:C:114:LYS:NZ	2:C:120:GLU:OE2	2.28	0.66
2:A:544:ILE:HG22	2:A:545:GLU:N	2.11	0.65
1:E:52:ASN:HD21	1:E:57:ARG:HB3	1.62	0.64
1:E:89:GLU:OE1	1:E:89:GLU:N	2.29	0.64
1:F:101:TRP:HH2	4:H:5:MAN:O4	1.80	0.64
2:B:566:THR:HG21	4:H:1:NAG:H82	1.80	0.62
2:C:163:ASP:OD1	2:C:543:TYR:OH	2.15	0.62
1:F:52:ASN:O	1:F:72:ARG:NH1	2.31	0.62
2:B:508:GLN:O	1:F:57:ARG:NH1	2.34	0.61
2:C:89:ARG:NH1	2:C:90:SER:O	2.33	0.61
2:C:115:LYS:HD2	2:C:119:SER:HB2	1.83	0.61
2:C:522:ASP:O	2:C:523:GLU:HG3	2.01	0.60
2:C:52:VAL:H	2:C:596:ARG:NH2	1.99	0.60
2:B:524:GLY:HA2	2:C:571:LEU:HD13	1.82	0.60
2:A:521:GLN:CD	2:A:544:ILE:HD11	2.23	0.60
2:A:76:ALA:O	2:A:85:ARG:NH2	2.34	0.59
1:E:52:ASN:ND2	1:E:57:ARG:HB3	2.17	0.59
2:B:163:ASP:OD1	2:B:543:TYR:OH	2.21	0.58
2:A:163:ASP:OD1	2:A:543:TYR:OH	2.22	0.58
1:F:91:THR:HG22	1:F:121:VAL:H	1.69	0.58
2:A:186:LEU:HD22	2:A:193:PHE:CE2	2.38	0.57
2:A:544:ILE:CG2	2:A:545:GLU:N	2.68	0.56
1:F:101:TRP:HH2	4:H:5:MAN:HO4	1.52	0.56
2:C:607:ASP:OD1	2:C:608:CYS:N	2.39	0.56
1:F:89:GLU:OE1	1:F:89:GLU:N	2.30	0.56
1:F:51:ILE:HG13	1:F:58:THR:HG22	1.88	0.56
2:A:550:ASN:OD1	2:A:551:GLN:N	2.38	0.56
2:B:607:ASP:OD1	2:B:608:CYS:N	2.39	0.55
1:F:6:GLU:OE1	1:F:6:GLU:N	2.38	0.55
2:A:607:ASP:OD1	2:A:608:CYS:N	2.40	0.55
2:B:187:PRO:HG2	2:B:193:PHE:CE2	2.41	0.55
6:E:201:MAN:HO6	5:I:4:MAN:HO6	1.40	0.54
2:C:64:ARG:HD2	2:C:186:LEU:HD11	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:201:MAN:H5	5:I:4:MAN:O5	2.08	0.54
2:C:50:LYS:HA	2:C:50:LYS:HE3	1.90	0.54
1:E:1:GLN:O	1:E:112:SER:OG	2.25	0.54
1:D:106:LEU:O	1:D:109:THR:HG22	2.08	0.54
2:B:189:ALA:O	2:B:191:LYS:NZ	2.40	0.54
2:C:45:VAL:HB	2:C:504:ILE:HD11	1.90	0.53
1:D:101:TRP:CZ2	3:G:3:MAN:H3	2.44	0.53
2:B:39:HIS:NE2	2:B:55:ASP:OD1	2.42	0.53
2:C:38:ILE:HG21	2:C:193:PHE:HE2	1.73	0.53
2:A:115:LYS:HD2	2:A:119:SER:HB2	1.90	0.52
1:F:12:VAL:HG11	1:F:86:LEU:HD13	1.91	0.52
1:D:99:ARG:HG3	1:D:110:TYR:HA	1.91	0.52
6:E:201:MAN:C5	5:I:4:MAN:C6	2.81	0.52
2:C:198:PRO:HD2	2:C:547:LEU:H	1.76	0.51
2:A:544:ILE:CG2	2:A:545:GLU:H	2.24	0.51
4:H:4:MAN:H2	4:H:5:MAN:C5	2.42	0.50
2:A:517:TYR:CZ	2:A:546:GLY:HA3	2.47	0.49
1:D:6:GLU:OE1	1:D:6:GLU:N	2.38	0.49
2:C:194:PHE:HA	2:C:517:TYR:CE1	2.47	0.49
1:D:52:ASN:OD1	1:D:57:ARG:HB3	2.12	0.48
2:A:599:GLY:HA2	2:C:597:TRP:HH2	1.78	0.48
1:D:91:THR:HG22	1:D:121:VAL:H	1.78	0.48
2:A:522:ASP:O	2:A:524:GLY:N	2.46	0.48
2:B:162:TYR:CE1	2:B:176:PHE:HB3	2.49	0.48
2:C:39:HIS:NE2	2:C:55:ASP:OD1	2.47	0.48
2:C:140:LYS:NZ	2:C:217:THR:HG23	2.28	0.48
2:B:49:ASP:OD1	2:B:50:LYS:N	2.46	0.48
2:B:180:VAL:HG23	2:B:562:ALA:HB1	1.95	0.48
1:E:104:ALA:CB	1:E:109:THR:HG21	2.44	0.47
2:C:53:CYS:O	2:C:56:LYS:NZ	2.43	0.47
2:C:89:ARG:HB3	2:C:153:PHE:CD2	2.49	0.47
1:D:22:CYS:HB3	1:D:79:VAL:HG22	1.97	0.47
4:H:4:MAN:H2	4:H:5:MAN:H5	1.97	0.47
2:A:53:CYS:SG	2:A:54:ARG:N	2.87	0.47
1:E:3:GLN:OE1	1:E:3:GLN:N	2.47	0.47
1:D:34:MET:HG3	1:D:79:VAL:HG11	1.96	0.46
2:C:563:ASN:OD1	5:I:1:NAG:N2	2.48	0.46
2:A:70:LEU:HG	2:A:179:GLY:HA2	1.96	0.46
2:C:63:LEU:HD23	2:C:185:ILE:HG12	1.97	0.46
1:F:34:MET:HG3	1:F:79:VAL:HG11	1.97	0.46
2:B:88:PHE:HB3	2:B:149:GLY:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:140:LYS:HD2	2:B:219:ARG:HH21	1.81	0.46
2:B:194:PHE:HA	2:B:517:TYR:CE2	2.51	0.46
2:B:198:PRO:HD2	2:B:547:LEU:H	1.81	0.46
2:A:562:ALA:HA	2:A:565:THR:HG22	1.97	0.45
2:C:182:ALA:HB2	2:C:562:ALA:HA	1.98	0.45
2:A:56:LYS:HA	2:A:56:LYS:HD3	1.80	0.45
1:D:101:TRP:O	1:D:102:SER:HB2	2.17	0.45
2:B:38:ILE:HB	2:B:187:PRO:HD3	1.98	0.45
1:E:102:SER:OG	1:E:103:ILE:N	2.50	0.44
2:A:97:VAL:O	2:A:166:ALA:N	2.45	0.44
1:E:100:PRO:O	1:E:101:TRP:C	2.59	0.44
1:F:52:ASN:OD1	1:F:57:ARG:HB3	2.16	0.44
2:A:599:GLY:HA2	2:C:597:TRP:CH2	2.51	0.44
2:B:127:ASP:N	2:B:127:ASP:OD1	2.50	0.44
2:A:162:TYR:CE2	2:A:176:PHE:HB3	2.53	0.44
2:A:596:ARG:CZ	2:A:596:ARG:HB2	2.48	0.44
2:B:56:LYS:HD3	2:B:56:LYS:HA	1.75	0.44
2:B:61:ASN:O	2:B:64:ARG:NH2	2.51	0.44
1:E:101:TRP:O	1:E:102:SER:HB2	2.18	0.44
1:E:115:GLN:H	1:E:115:GLN:CD	2.26	0.43
2:C:127:ASP:OD1	2:C:128:GLY:N	2.50	0.43
2:B:529:LEU:HB3	2:B:532:ILE:HD12	2.00	0.43
2:B:527:ILE:HG23	2:B:528:GLY:N	2.33	0.43
1:E:34:MET:HG3	1:E:79:VAL:HG11	2.01	0.43
2:A:97:VAL:HG22	2:A:573:LEU:HD21	2.01	0.43
4:H:4:MAN:C2	4:H:5:MAN:H5	2.49	0.43
2:A:96:VAL:CG2	2:A:580:ARG:HG2	2.49	0.43
2:A:88:PHE:HB3	2:A:149:GLY:H	1.83	0.42
2:C:122:LEU:HB2	2:C:172:ARG:HB2	2.01	0.42
2:A:513:PRO:C	2:A:555:ILE:HD11	2.44	0.42
2:A:591:ASP:O	2:A:595:GLN:HG2	2.19	0.42
2:C:69:ASN:HA	2:C:179:GLY:HA3	2.01	0.42
2:C:140:LYS:HZ2	2:C:217:THR:HG23	1.84	0.42
2:B:597:TRP:HH2	2:C:599:GLY:HA2	1.85	0.42
1:D:30:SER:O	1:D:53:TYR:HB2	2.19	0.42
2:A:127:ASP:OD1	2:A:127:ASP:N	2.53	0.42
2:C:601:CYS:HB2	2:C:608:CYS:HB2	1.39	0.42
1:D:99:ARG:HG2	1:D:100:PRO:HD2	2.00	0.42
1:E:106:LEU:O	1:E:109:THR:HG22	2.19	0.42
1:D:91:THR:HG22	1:D:120:THR:HA	2.01	0.41
2:A:39:HIS:O	2:A:42:THR:OG1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:64:ARG:HD3	2:A:64:ARG:HA	1.87	0.41
2:B:601:CYS:HB2	2:B:607:ASP:OD1	2.20	0.41
2:C:162:TYR:CE2	2:C:176:PHE:HB3	2.55	0.41
2:A:96:VAL:HG12	2:A:167:SER:HB2	2.01	0.41
2:A:194:PHE:HA	2:A:517:TYR:CZ	2.55	0.41
1:D:102:SER:OG	1:D:103:ILE:N	2.54	0.41
2:B:597:TRP:CH2	2:C:599:GLY:HA2	2.56	0.41
2:A:544:ILE:HG22	2:A:545:GLU:H	1.83	0.41
1:D:100:PRO:HG2	2:C:528:GLY:O	2.20	0.41
2:C:521:GLN:HB2	2:C:544:ILE:HG13	2.01	0.41
1:D:91:THR:HG23	1:D:121:VAL:HG12	2.03	0.40
2:B:106:GLU:OE1	2:B:136:ARG:NH1	2.54	0.40
2:C:97:VAL:HG23	2:C:573:LEU:HD11	2.04	0.40
1:E:89:GLU:H	1:E:89:GLU:CD	2.26	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	D	121/144 (84%)	116 (96%)	3 (2%)	2 (2%)	7	30
1	E	121/144 (84%)	116 (96%)	3 (2%)	2 (2%)	7	30
1	F	121/144 (84%)	116 (96%)	3 (2%)	2 (2%)	7	30
2	A	287/706 (41%)	272 (95%)	14 (5%)	1 (0%)	36	67
2	B	287/706 (41%)	274 (96%)	12 (4%)	1 (0%)	36	67
2	C	287/706 (41%)	270 (94%)	16 (6%)	1 (0%)	36	67
All	All	1224/2550 (48%)	1164 (95%)	51 (4%)	9 (1%)	20	49

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	102	SER
2	A	523	GLU
1	E	102	SER
1	D	2	VAL
1	F	56	GLU
1	F	101	TRP
2	B	523	GLU
2	C	523	GLU
1	E	2	VAL

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	D	96/112 (86%)	93 (97%)	3 (3%)	35	64
1	E	96/112 (86%)	94 (98%)	2 (2%)	47	71
1	F	96/112 (86%)	95 (99%)	1 (1%)	68	79
2	A	240/598 (40%)	240 (100%)	0	100	100
2	B	240/598 (40%)	240 (100%)	0	100	100
2	C	240/598 (40%)	239 (100%)	1 (0%)	84	86
All	All	1008/2130 (47%)	1001 (99%)	7 (1%)	73	82

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	99	ARG
1	D	101	TRP
1	D	103	ILE
2	C	544	ILE
1	E	101	TRP
1	E	103	ILE
1	F	103	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	13	GLN
1	D	118	GLN
2	A	62	GLN
2	A	188	GLN
2	A	514	ASN
2	A	567	GLN
2	B	62	GLN
2	B	73	ASN
2	B	188	GLN
2	B	197	HIS
2	B	508	GLN
2	B	567	GLN
2	C	62	GLN
2	C	508	GLN
2	C	560	GLN
1	E	13	GLN
1	E	52	ASN
1	E	82	GLN
1	F	82	GLN
1	F	84	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](#)

20 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	3,2	14,14,15	0.40	0	17,19,21	0.45	0
3	NAG	G	2	3	14,14,15	0.41	0	17,19,21	0.63	0
3	MAN	G	3	3	11,11,12	0.44	0	15,15,17	0.47	0
3	MAN	G	4	3	11,11,12	0.42	0	15,15,17	1.07	2 (13%)
3	MAN	G	5	3	11,11,12	0.21	0	15,15,17	0.66	1 (6%)
3	MAN	G	6	3	11,11,12	0.29	0	15,15,17	0.36	0
3	MAN	G	7	3	11,11,12	0.34	0	15,15,17	0.48	0
4	NAG	H	1	4,2	14,14,15	0.40	0	17,19,21	0.60	0
4	NAG	H	2	4	14,14,15	0.41	0	17,19,21	0.45	0
4	BMA	H	3	4	11,11,12	0.40	0	15,15,17	0.51	0
4	MAN	H	4	4	11,11,12	0.51	0	15,15,17	1.14	2 (13%)
4	MAN	H	5	4	11,11,12	0.33	0	15,15,17	0.56	0
4	BMA	H	6	4	11,11,12	0.22	0	15,15,17	0.47	0
4	MAN	H	7	4	11,11,12	0.29	0	15,15,17	0.42	0
5	NAG	I	1	5,2	14,14,15	0.45	0	17,19,21	0.90	1 (5%)
5	NAG	I	2	5	14,14,15	0.41	0	17,19,21	0.37	0
5	BMA	I	3	5	11,11,12	0.36	0	15,15,17	0.47	0
5	MAN	I	4	5	11,11,12	0.30	0	15,15,17	0.48	0
5	MAN	I	5	5	11,11,12	0.23	0	15,15,17	0.42	0
5	MAN	I	6	5	11,11,12	0.28	0	15,15,17	0.46	0

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	3,2	-	4/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	MAN	G	3	3	-	1/2/19/22	0/1/1/1
3	MAN	G	4	3	-	0/2/19/22	0/1/1/1
3	MAN	G	5	3	-	0/2/19/22	0/1/1/1
3	MAN	G	6	3	-	0/2/19/22	1/1/1/1
3	MAN	G	7	3	-	0/2/19/22	0/1/1/1
4	NAG	H	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	BMA	H	3	4	-	2/2/19/22	0/1/1/1
4	MAN	H	4	4	-	2/2/19/22	0/1/1/1
4	MAN	H	5	4	-	1/2/19/22	0/1/1/1
4	BMA	H	6	4	-	0/2/19/22	0/1/1/1
4	MAN	H	7	4	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	I	1	5,2	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	0/6/23/26	0/1/1/1
5	BMA	I	3	5	-	2/2/19/22	0/1/1/1
5	MAN	I	4	5	-	2/2/19/22	0/1/1/1
5	MAN	I	5	5	-	0/2/19/22	0/1/1/1
5	MAN	I	6	5	-	0/2/19/22	1/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	4	MAN	C1-O5-C5	2.90	116.07	112.19
4	H	4	MAN	C3-C4-C5	2.44	114.66	110.23
5	I	1	NAG	C1-C2-N2	2.27	114.01	110.43
3	G	4	MAN	C1-C2-C3	2.18	112.82	109.64
4	H	4	MAN	O5-C1-C2	-2.13	105.72	110.79
3	G	5	MAN	C1-C2-C3	2.05	112.63	109.64

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	3	BMA	C4-C5-C6-O6
4	H	3	BMA	O5-C5-C6-O6
4	H	4	MAN	O5-C5-C6-O6
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
5	I	4	MAN	C4-C5-C6-O6
5	I	4	MAN	O5-C5-C6-O6
5	I	3	BMA	C4-C5-C6-O6
4	H	2	NAG	C8-C7-N2-C2
3	G	3	MAN	O5-C5-C6-O6
4	H	5	MAN	O5-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
4	H	4	MAN	C4-C5-C6-O6
4	H	2	NAG	O7-C7-N2-C2
5	I	3	BMA	O5-C5-C6-O6

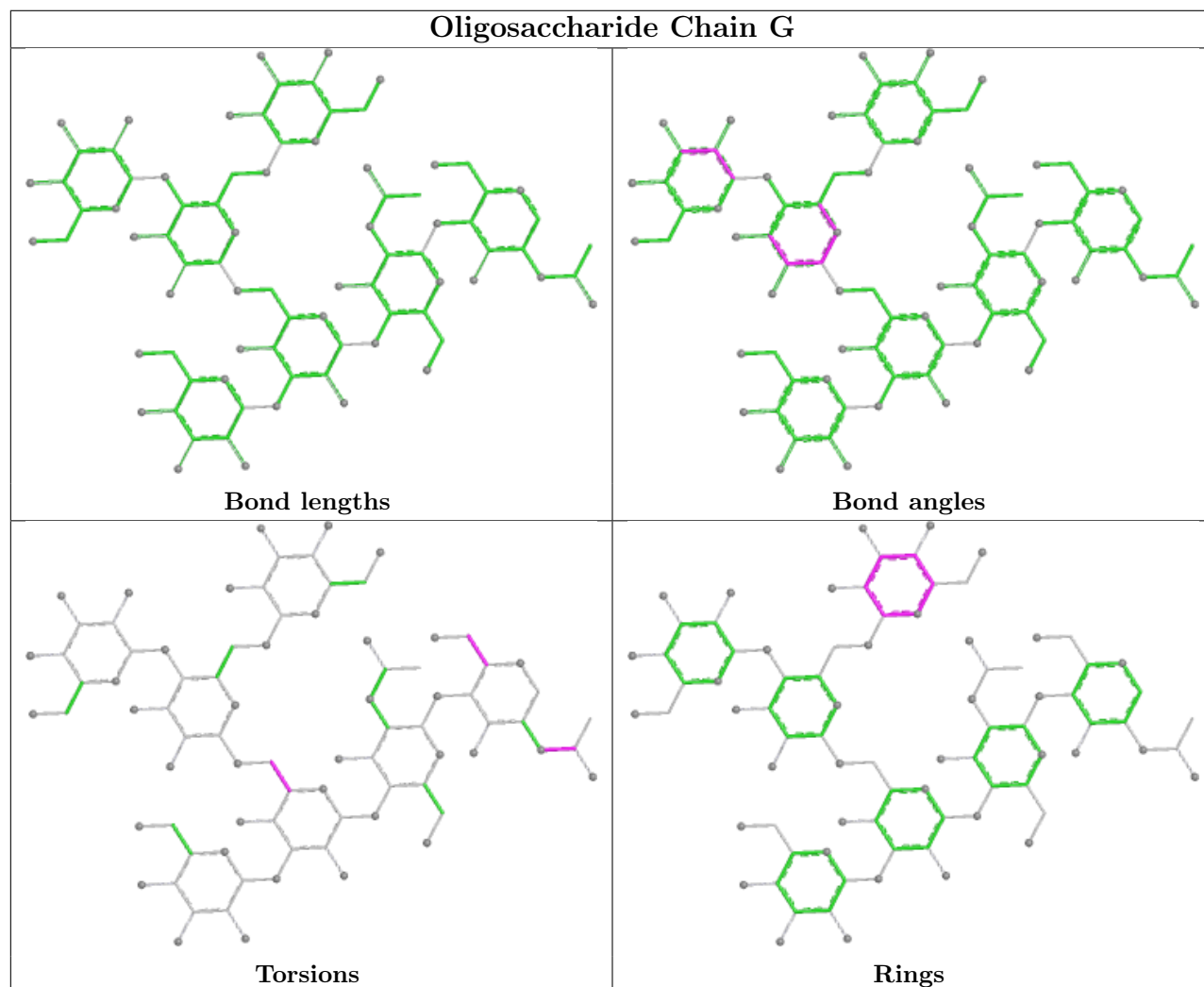
All (2) ring outliers are listed below:

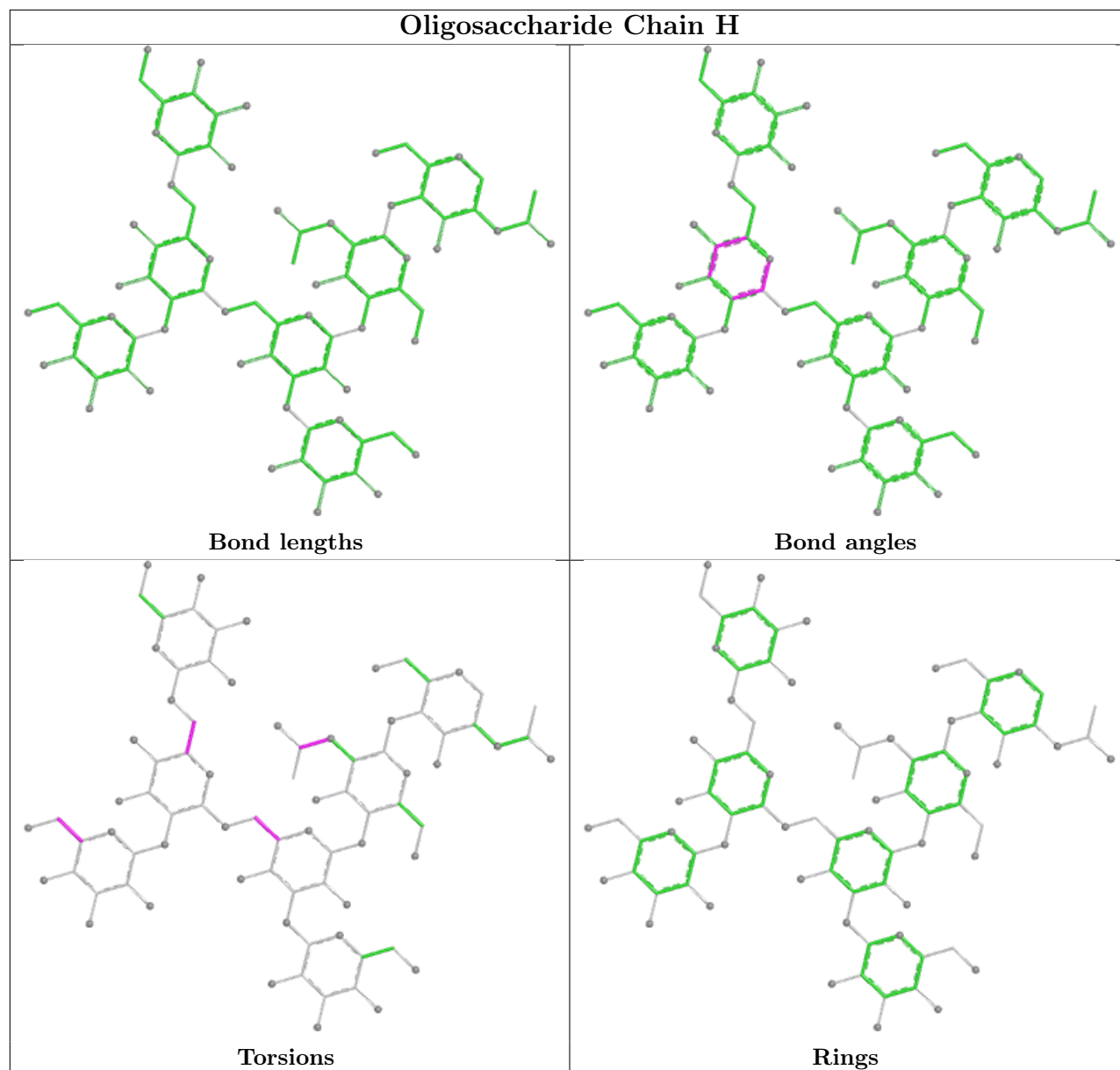
Mol	Chain	Res	Type	Atoms
3	G	6	MAN	C1-C2-C3-C4-C5-O5
5	I	6	MAN	C1-C2-C3-C4-C5-O5

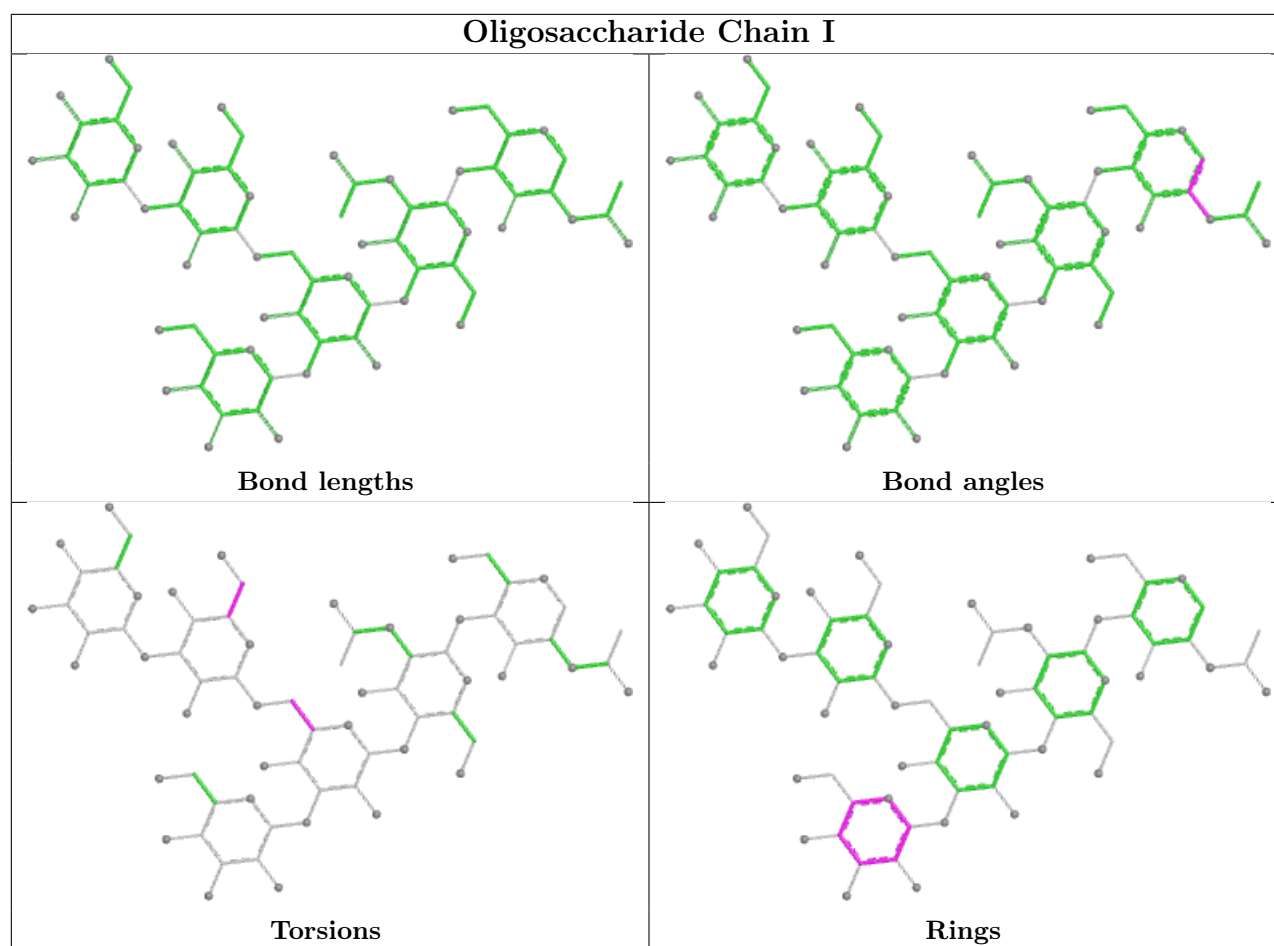
6 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	3	MAN	1	0
4	H	4	MAN	3	0
4	H	5	MAN	5	0
5	I	4	MAN	12	0
4	H	1	NAG	1	0
5	I	1	NAG	1	0

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 [i](#)

1 ligand is 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	$\# Z > 2$
6	MAN	E	201	-	11,11,12	0.23	0	15,15,17	0.50	0

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
6	MAN	E	201	-	-	2/2/18/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	E	201	MAN	C4-C5-C6-O6
6	E	201	MAN	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	201	MAN	12	0

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.

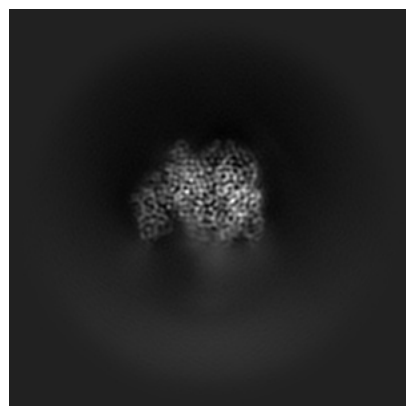
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44873. These allow visual inspection of the internal detail of the map and identification of artifacts.

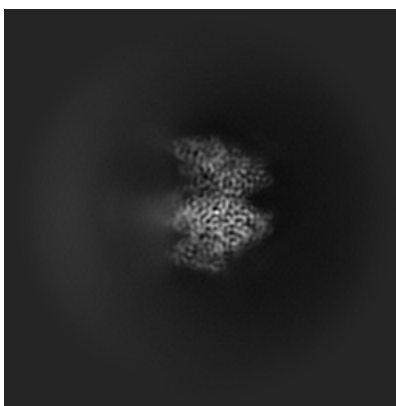
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

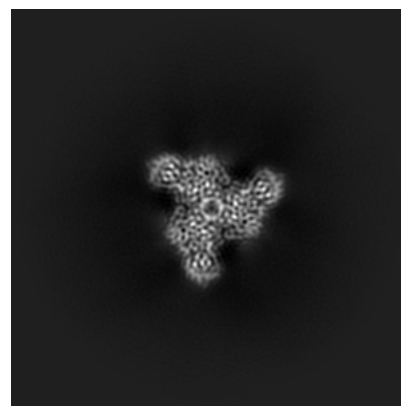
6.1.1 Primary map



X

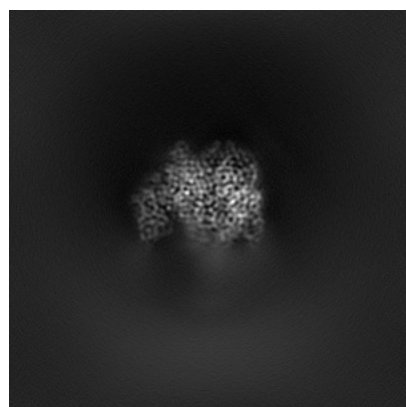


Y

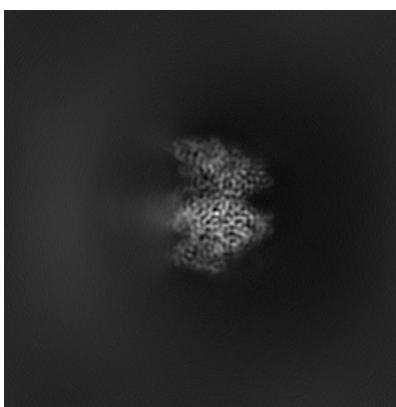


Z

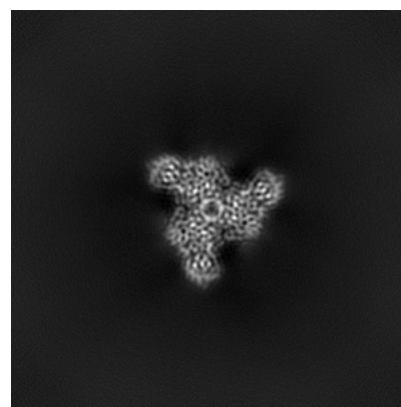
6.1.2 Raw map



X



Y



Z

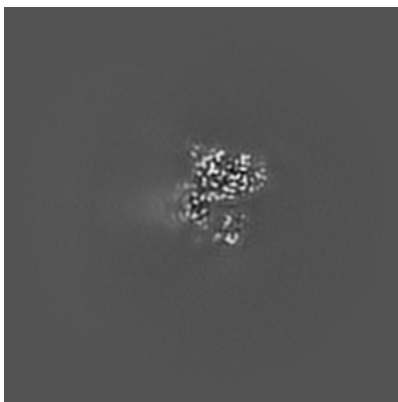
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

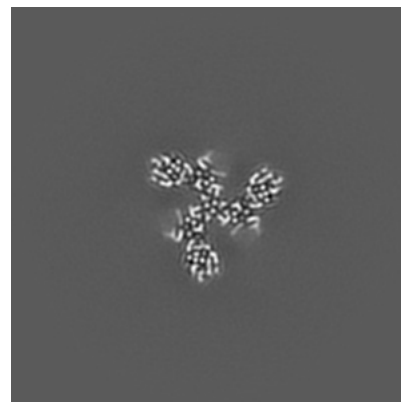
6.2.1 Primary map



X Index: 160

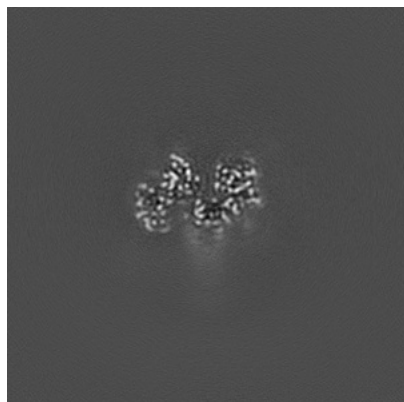


Y Index: 160

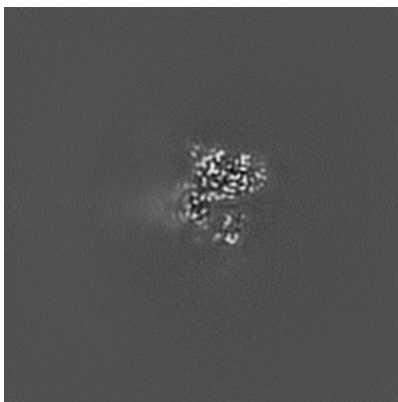


Z Index: 160

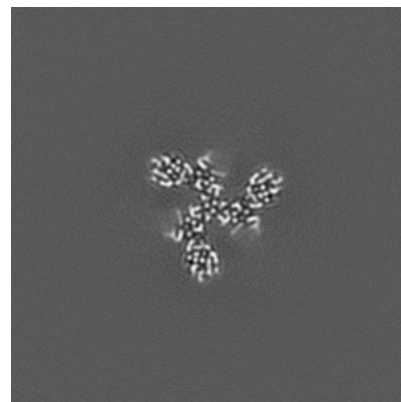
6.2.2 Raw map



X Index: 160



Y Index: 160

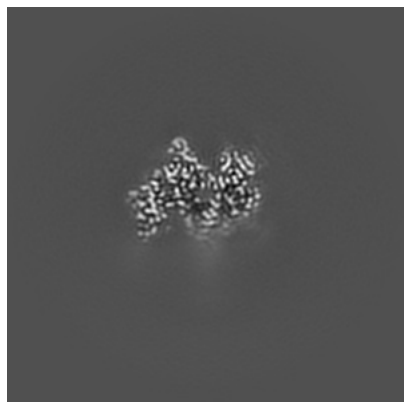


Z Index: 160

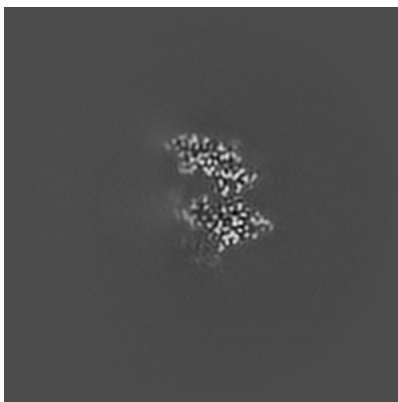
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

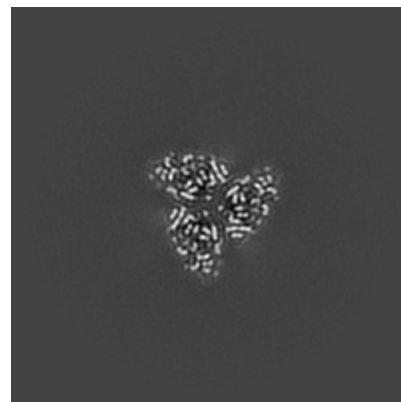
6.3.1 Primary map



X Index: 153

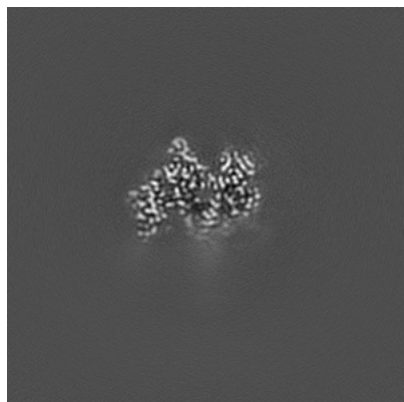


Y Index: 174

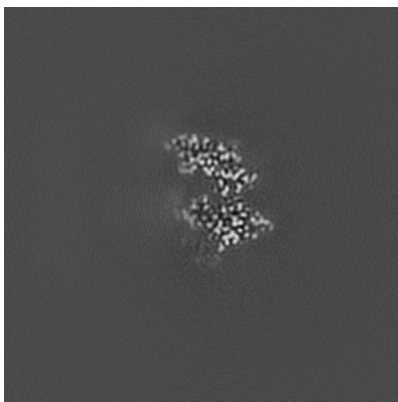


Z Index: 173

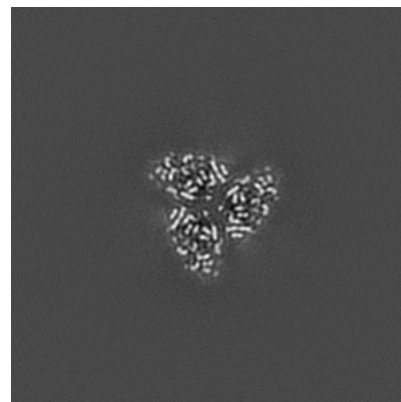
6.3.2 Raw map



X Index: 153



Y Index: 174

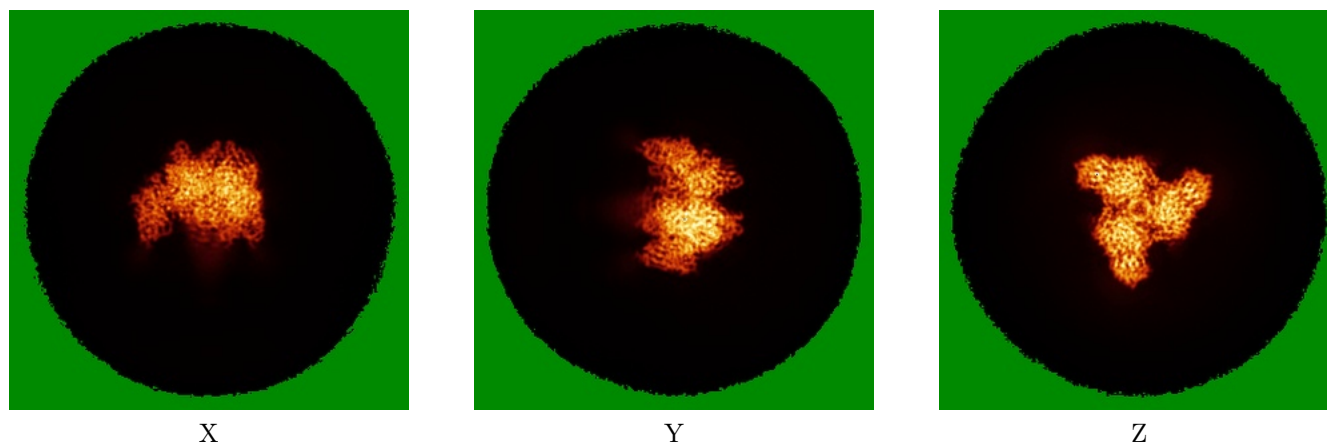


Z Index: 173

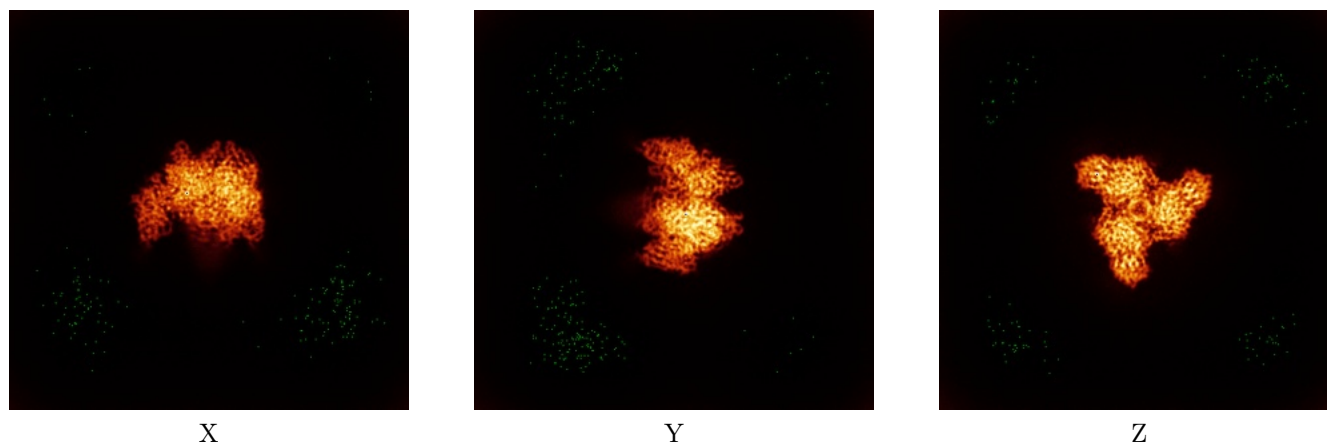
The images above show the largest variance slices of the map in three orthogonal directions.

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.

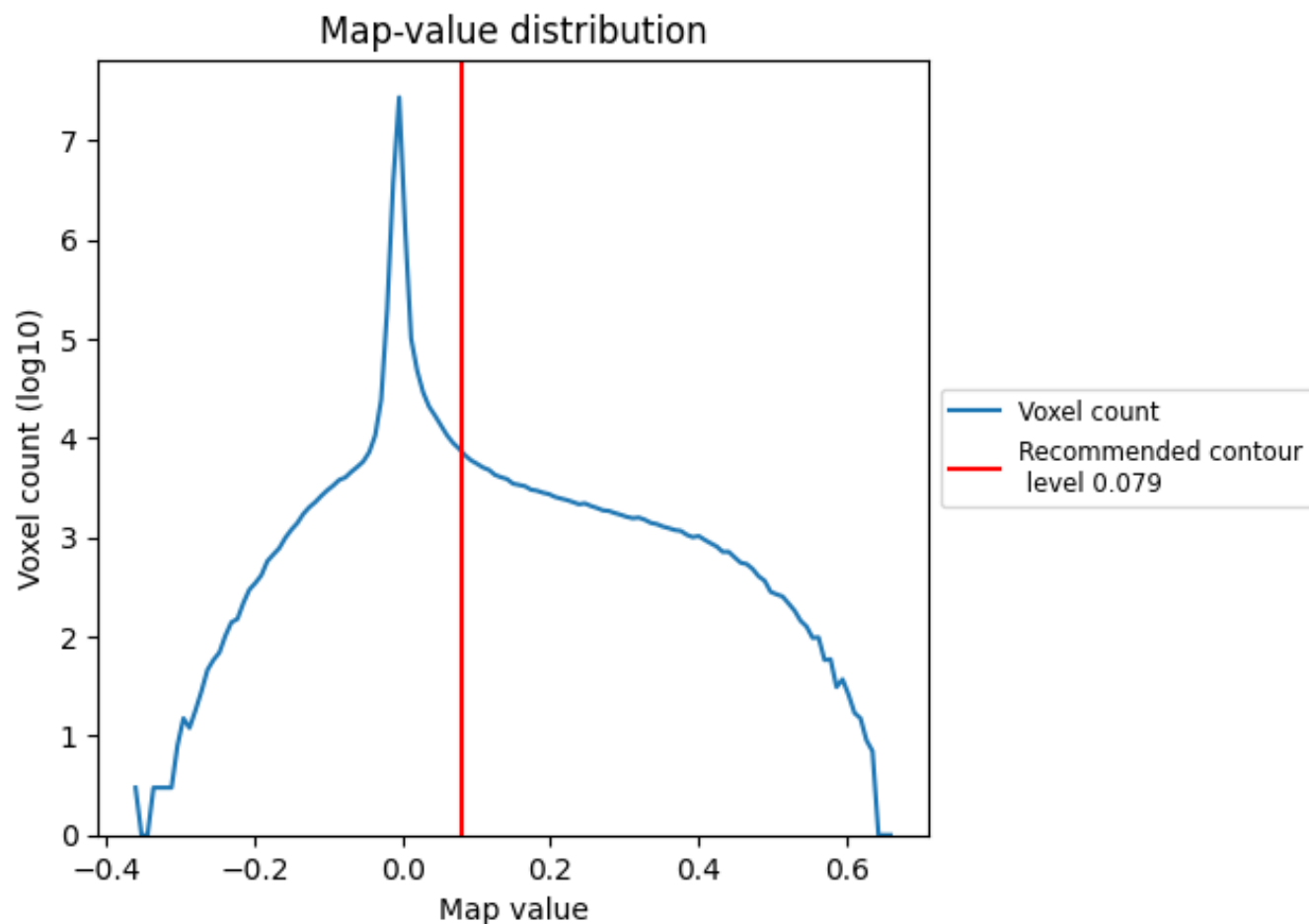
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

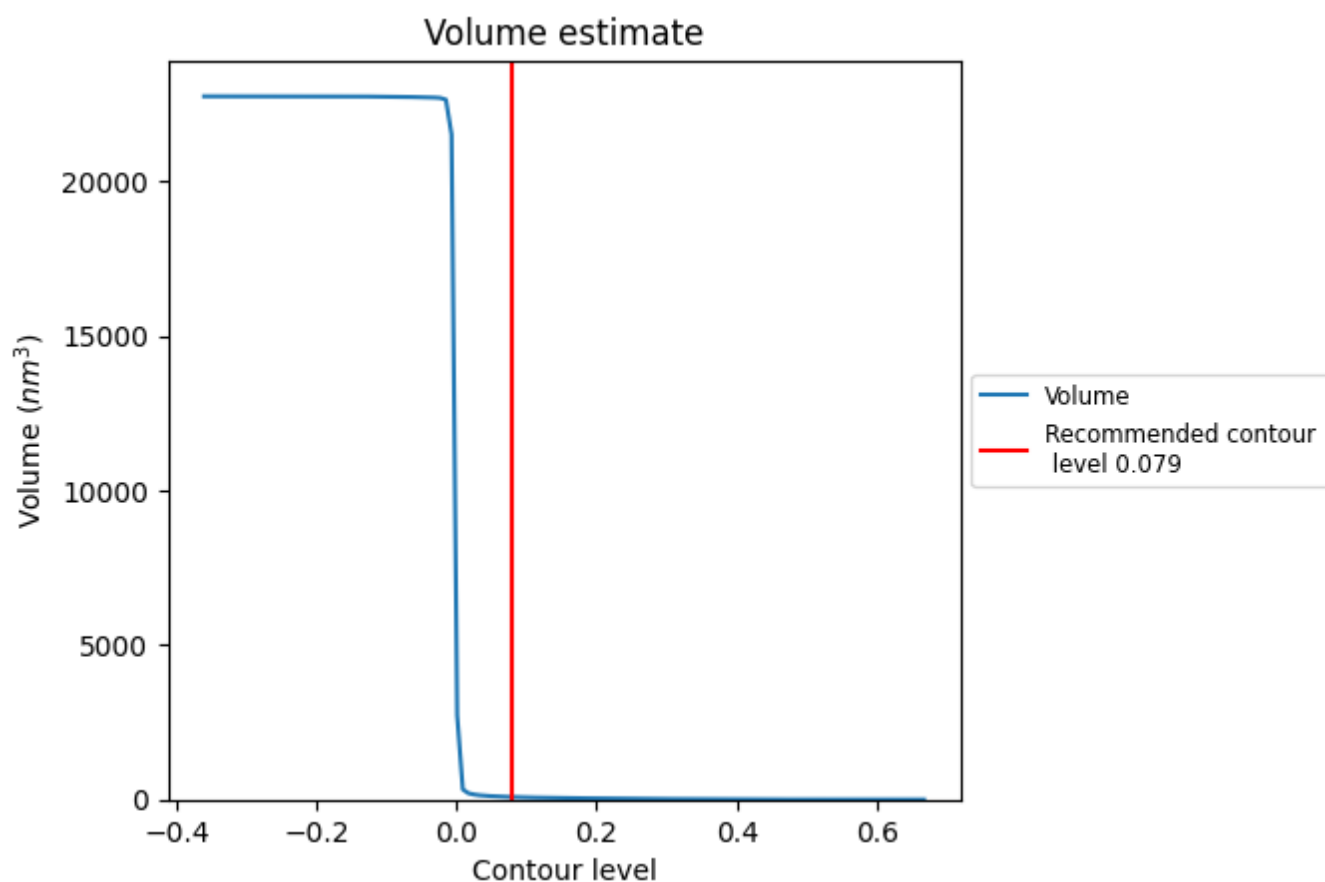
This section contains the results of statistical analysis of the map.

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.

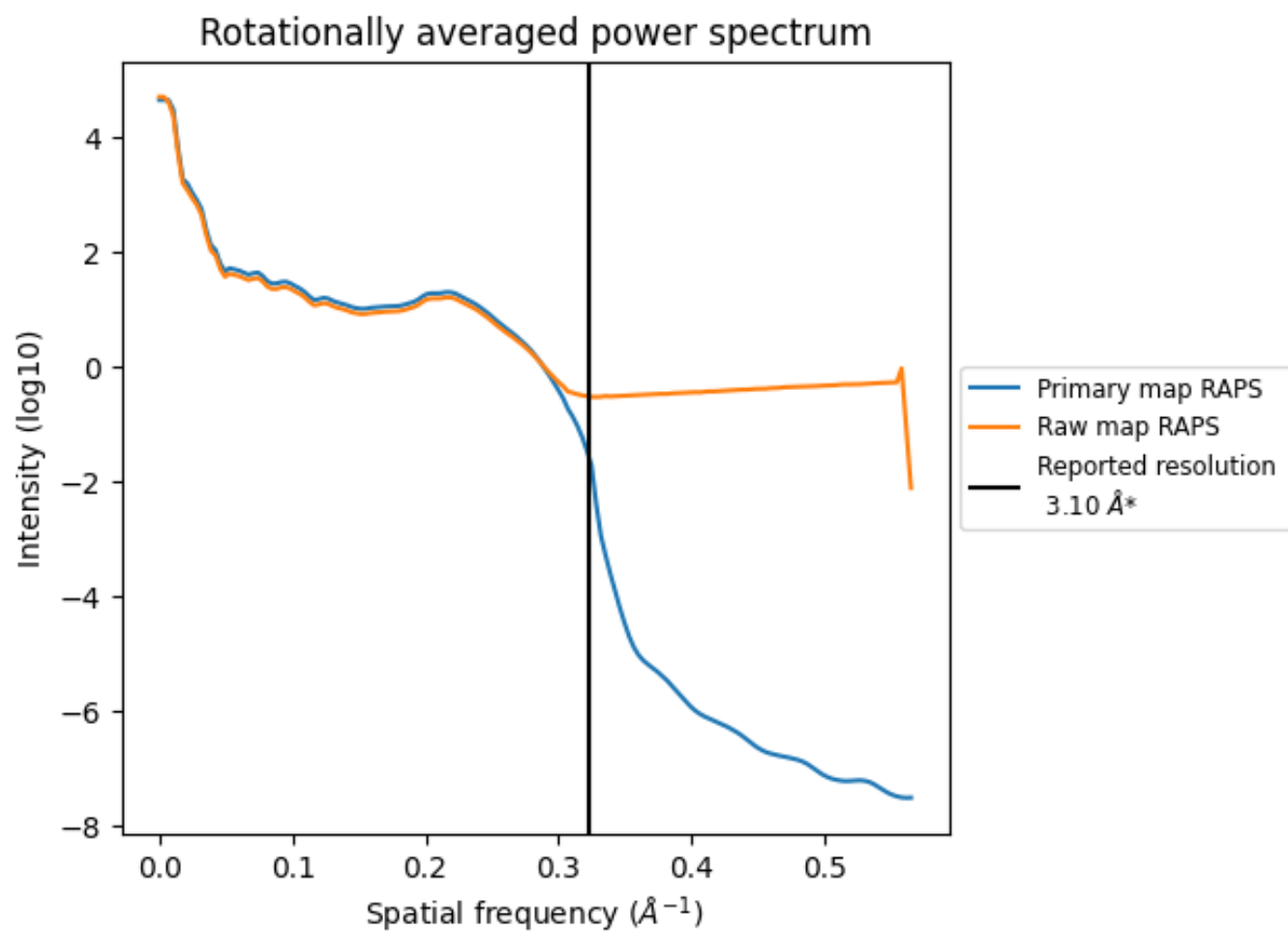
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 82 nm³; this corresponds to an approximate mass of 74 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 ⓘ

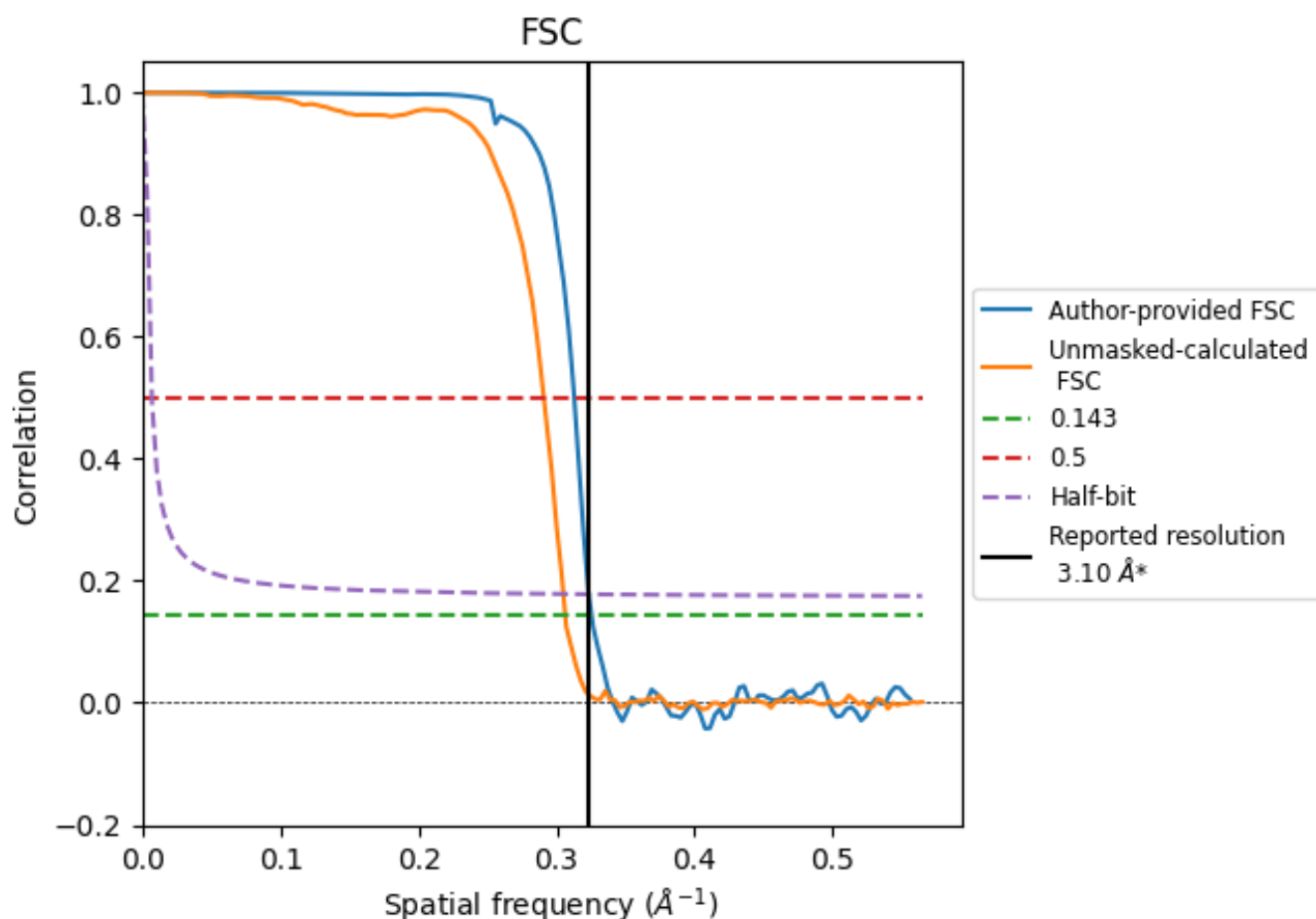


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 $0.323 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.07	3.20	3.09
Unmasked-calculated*	3.26	3.44	3.28

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

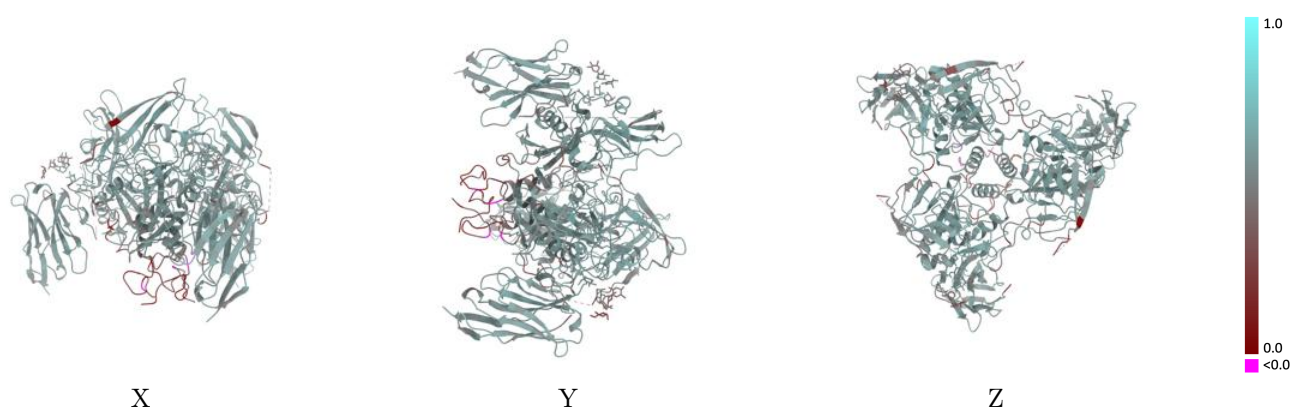
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-44873 and PDB model 9BSV. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)

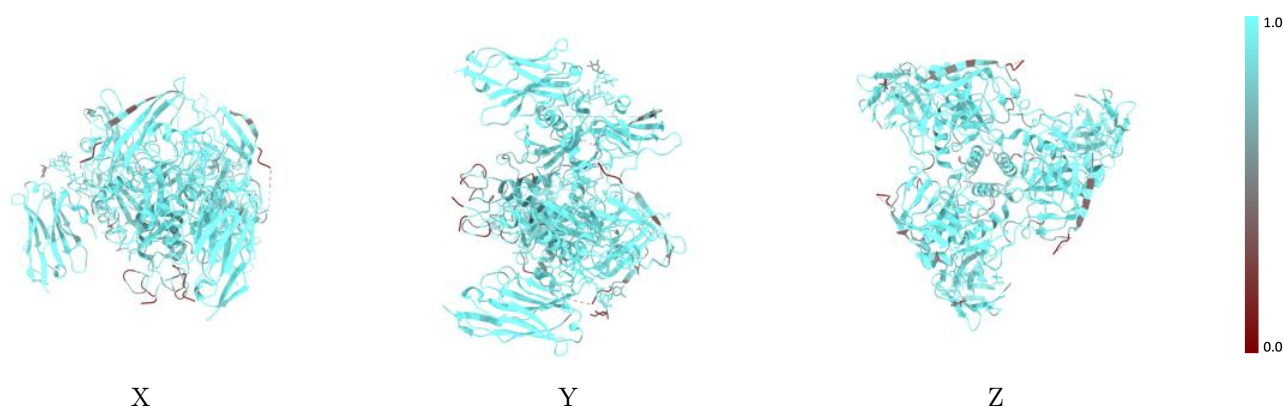
This section was not generated.

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



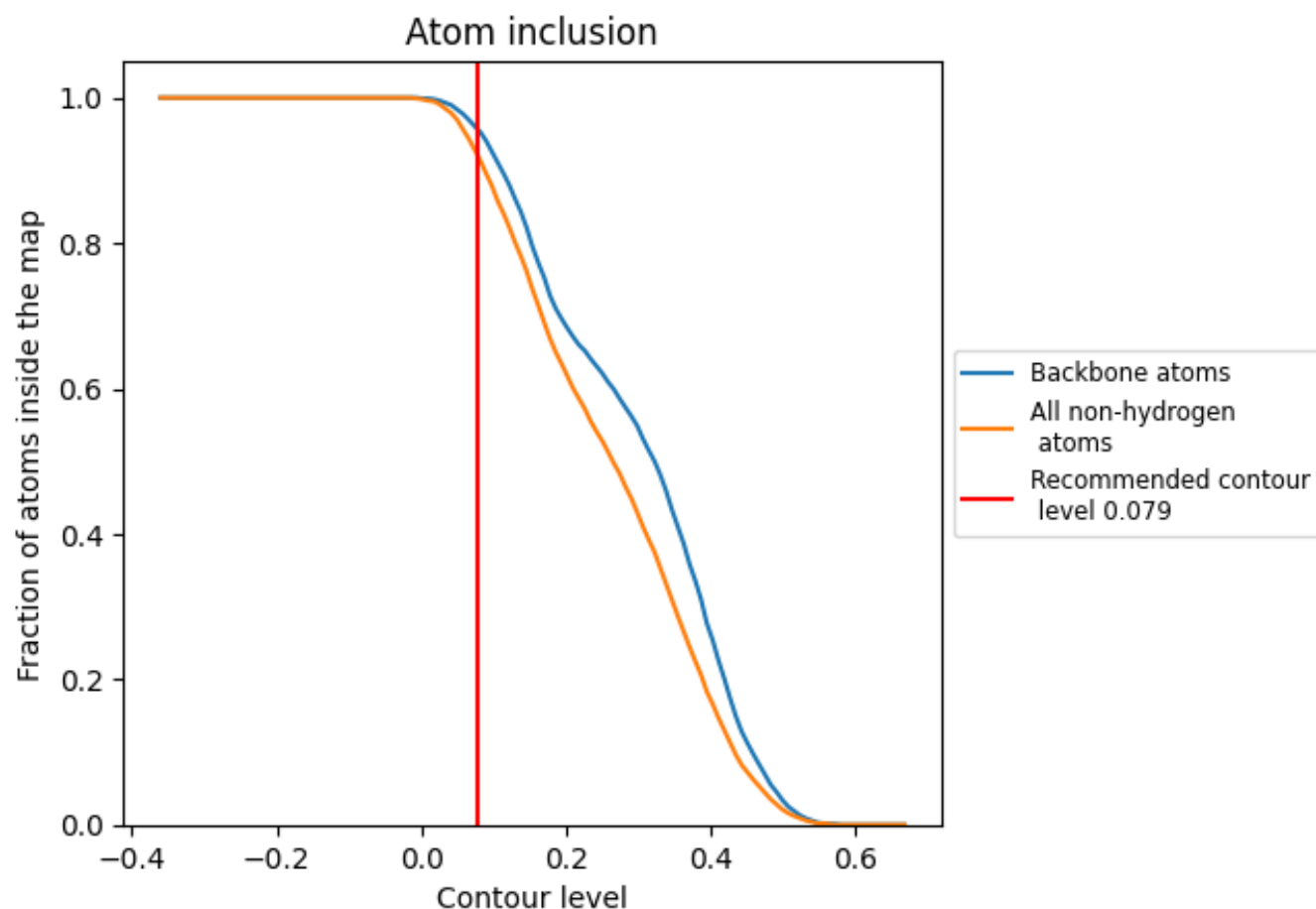
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



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 (0.079).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 92% 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 (0.079) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9180	0.5360
A	0.9000	0.5270
B	0.9020	0.5320
C	0.9060	0.5280
D	0.9590	0.5570
E	0.9550	0.5580
F	0.9610	0.5580
G	0.8680	0.4750
H	0.8310	0.4470
I	0.9860	0.5380

1.0

0.0

<0.0