



Full wwPDB EM Validation Report ⓘ

Mar 14, 2026 – 07:01 AM UTC

PDB ID : 8PC0 / pdb_00008pc0
EMDB ID : EMD-17592
Title : Sub-tomogram average of the open conformation of the Nap adhesion complex from the human pathogen *Mycoplasma genitalium*.
Authors : Sprankel, L.; Scheffer, M.P.; Frangakis, A.F.
Deposited on : 2023-06-09
Resolution : 17.00 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 EMD 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

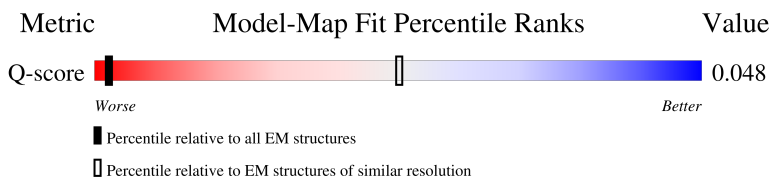
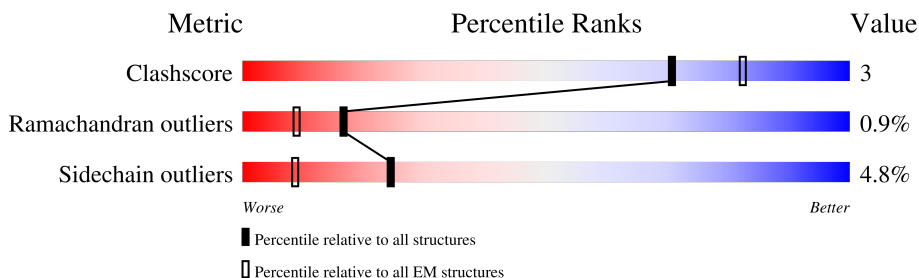
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 17.00 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	38 (16.50 - 17.50)

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	1053	
2	C	1444	
3	B	3	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 33611 atoms, of which 16590 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.

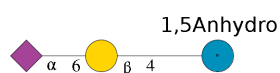
- Molecule 1 is a protein called Mgp-operon protein 3.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	887	Total	C	H	N	O	S	0	0
			13436	4281	6611	1138	1400	6		

- Molecule 2 is a protein called Adhesin P1.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	C	1285	Total	C	H	N	O	S	0	0
			19793	6365	9744	1699	1971	14		

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-1,5-anhydro-D-glucitol.



Mol	Chain	Residues	Atoms						AltConf	Trace
3	B	3	Total	C	H	N	O		0	0
			79	23	37	1	18			

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	K	0
			1	1	

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	O	P	0
			5	4	1	

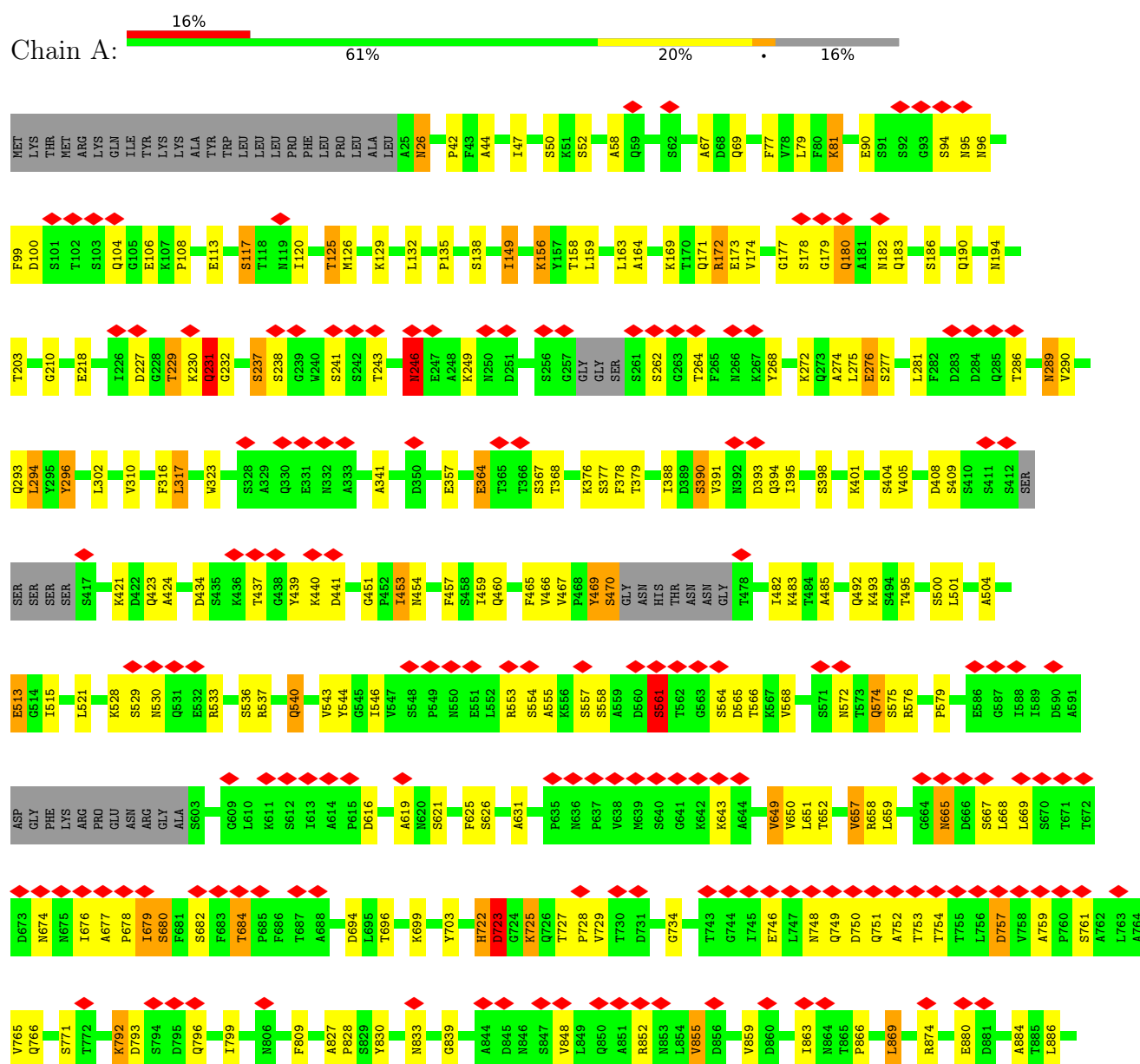
- Molecule 6 is water.

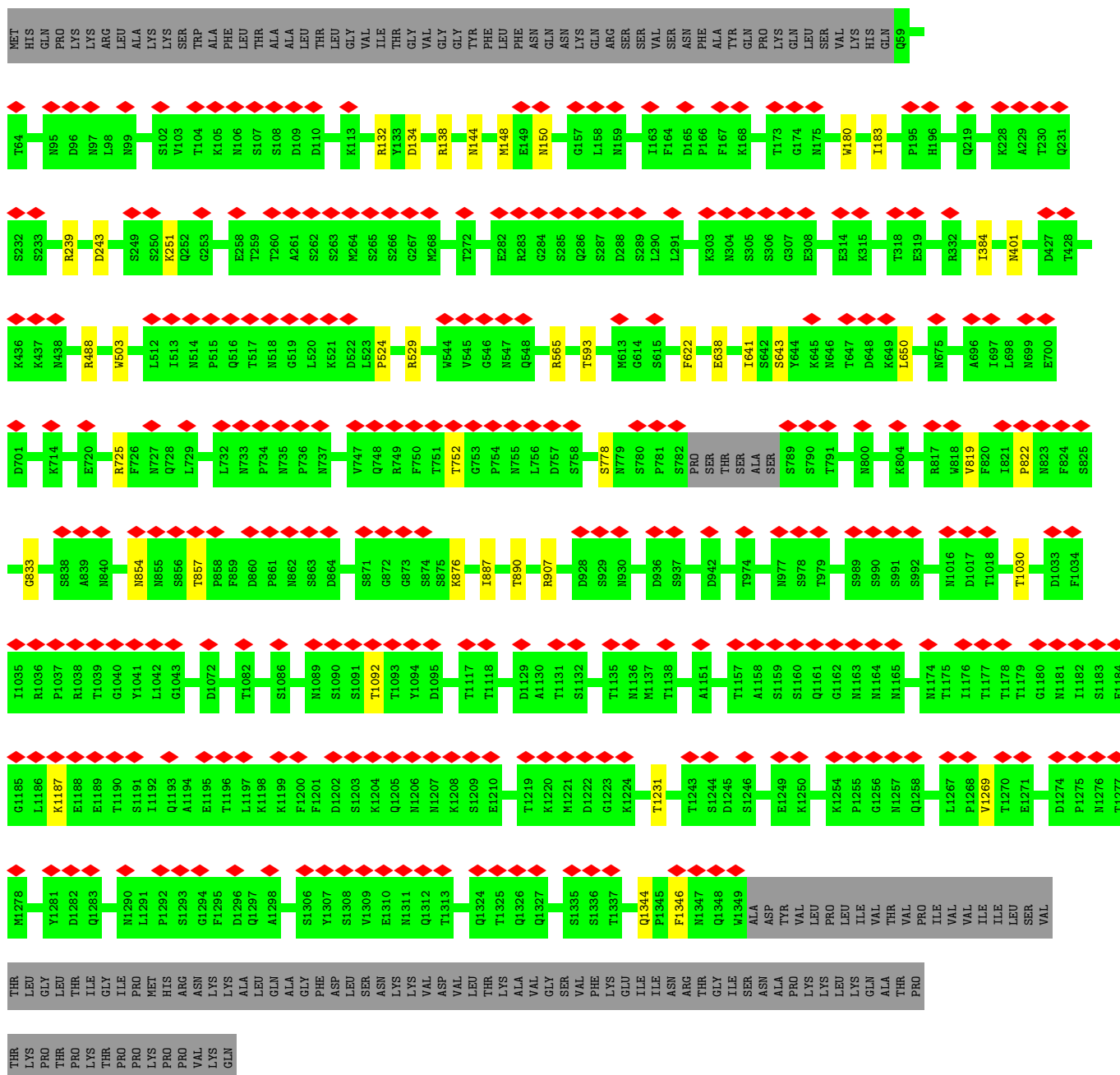
Mol	Chain	Residues	Atoms			AltConf
6	A	99	Total	H	O	0
			297	198	99	

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: Mgp-operon protein 3





- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-1,5-anhydro-D-glucitol

Chain B: 

AS01
GAL2
SIA3

4 Experimental information

Property	Value	Source
EM reconstruction method	SUBTOMOGRAM AVERAGING	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of subtomograms used	9689	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 ($e^-/\text{\AA}^2$)	120	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	4000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.001	Depositor
Minimum map value	-0.001	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.000229	Depositor
Map size (Å)	332.8, 332.8, 332.8	wwPDB
Map dimensions	64, 64, 64	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	5.2, 5.2, 5.2	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: K, GAL, PO4, ASO, SIA

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	1.17	11/6964 (0.2%)	1.53	49/9472 (0.5%)
2	C	0.89	0/10303	1.36	14/14029 (0.1%)
All	All	1.01	11/17267 (0.1%)	1.43	63/23501 (0.3%)

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	A	0	5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	927	SER	C-O	-8.68	1.12	1.23
1	A	401	LYS	C-O	6.94	1.32	1.24
1	A	923	ILE	C-O	6.89	1.32	1.24
1	A	649	VAL	C-O	6.15	1.30	1.23
1	A	920	PRO	C-O	-6.00	1.16	1.23
1	A	504	ALA	C-O	6.00	1.32	1.23
1	A	316	PHE	C-O	-5.89	1.17	1.24
1	A	404	SER	C-O	5.40	1.30	1.24
1	A	880	GLU	C-O	5.35	1.30	1.23
1	A	246	ASN	C-O	5.32	1.30	1.24
1	A	828	PRO	C-O	-5.16	1.17	1.23

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	827	ALA	O-C-N	-6.88	117.31	121.71
1	A	231	GLN	CA-C-N	-6.66	113.97	121.83
1	A	231	GLN	C-N-CA	-6.66	113.97	121.83
1	A	408	ASP	CA-CB-CG	6.61	119.21	112.60
1	A	95	ASN	CA-C-N	6.49	128.88	120.44
1	A	95	ASN	C-N-CA	6.49	128.88	120.44
1	A	513	GLU	CA-C-N	6.38	128.18	120.13
1	A	513	GLU	C-N-CA	6.38	128.18	120.13
1	A	723	ASP	CA-CB-CG	6.38	118.98	112.60
1	A	894	THR	CA-CB-OG1	-6.30	100.14	109.60
1	A	903	GLN	N-CA-CB	6.18	119.39	110.06
1	A	26	ASN	CA-CB-CG	6.12	118.72	112.60
1	A	723	ASP	CB-CA-C	-6.11	98.27	110.42
1	A	616	ASP	CA-CB-CG	6.03	118.63	112.60
1	A	616	ASP	CB-CA-C	5.94	120.27	111.06
2	C	138	ARG	NE-CZ-NH2	5.91	124.52	119.20
1	A	274	ALA	CA-C-N	5.90	128.44	120.65
1	A	274	ALA	C-N-CA	5.90	128.44	120.65
1	A	100	ASP	CA-C-N	5.89	128.09	120.44
1	A	100	ASP	C-N-CA	5.89	128.09	120.44
1	A	703	TYR	CA-C-O	-5.81	114.43	120.70
1	A	903	GLN	CB-CA-C	-5.77	99.77	109.53
1	A	42	PRO	CB-CA-C	-5.71	103.28	111.62
1	A	553	ARG	CA-C-O	-5.70	114.38	120.42
1	A	757	ASP	CA-CB-CG	5.61	118.21	112.60
1	A	722	HIS	CA-C-N	5.61	132.25	121.54
1	A	722	HIS	C-N-CA	5.61	132.25	121.54
2	C	384	ILE	N-CA-C	5.58	116.22	110.36
2	C	132	ARG	NE-CZ-NH2	5.58	124.22	119.20
2	C	907	ARG	NE-CZ-NH2	5.55	124.19	119.20
2	C	529	ARG	NE-CZ-NH2	5.48	124.13	119.20
1	A	809	PHE	CA-C-O	-5.43	115.08	121.16
1	A	923	ILE	CA-C-N	-5.38	115.21	122.63
1	A	923	ILE	C-N-CA	-5.38	115.21	122.63
1	A	238	SER	CA-C-N	5.31	125.87	119.98
1	A	238	SER	C-N-CA	5.31	125.87	119.98
2	C	524	PRO	N-CA-CB	5.30	106.29	102.79
1	A	316	PHE	CA-C-O	-5.29	114.15	120.56
2	C	134	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	754	THR	CA-C-N	5.28	129.04	121.50
1	A	754	THR	C-N-CA	5.28	129.04	121.50
2	C	150	ASN	CB-CA-C	5.25	114.93	111.20
1	A	117	SER	CA-C-O	-5.22	114.45	120.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	388	ILE	CA-C-O	-5.19	116.08	120.96
1	A	621	SER	CA-C-O	-5.19	115.30	121.32
1	A	729	VAL	CA-C-O	-5.19	114.93	120.59
1	A	684	THR	CB-CA-C	5.18	117.08	110.13
1	A	171	GLN	N-CA-C	-5.14	105.68	111.28
1	A	276	GLU	CA-C-N	5.13	127.47	120.54
1	A	276	GLU	C-N-CA	5.13	127.47	120.54
1	A	699	LYS	N-CA-CB	5.13	117.42	110.59
2	C	488	ARG	NE-CZ-NH2	5.12	123.81	119.20
1	A	839	GLY	CA-C-O	-5.12	115.92	122.01
1	A	874	ARG	CB-CA-C	5.11	119.06	109.71
1	A	289	ASN	CA-C-N	5.11	126.99	120.56
1	A	289	ASN	C-N-CA	5.11	126.99	120.56
2	C	854	ASN	CA-CB-CG	5.11	117.71	112.60
2	C	1269	VAL	CA-C-N	5.11	127.63	120.28
2	C	1269	VAL	C-N-CA	5.11	127.63	120.28
2	C	239	ARG	NE-CZ-NH2	5.10	123.79	119.20
1	A	296	TYR	CA-C-O	-5.09	115.15	120.55
1	A	833	ASN	N-CA-C	-5.09	106.92	112.72
2	C	565	ARG	NE-CZ-NH2	5.03	123.72	119.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	135	PRO	Peptide
1	A	178	SER	Peptide
1	A	179	GLY	Peptide
1	A	243	THR	Peptide
1	A	722	HIS	Peptide

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	6825	6611	6606	112	0
2	C	10049	9744	9727	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	42	37	37	1	0
4	A	1	0	0	0	0
5	A	5	0	0	0	0
6	A	99	198	0	2	0
All	All	17021	16590	16370	115	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 3.

All (115) 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:A:281:LEU:H	1:A:293:GLN:HE22	1.21	0.88
1:A:682:SER:OG	1:A:761:SER:HB3	1.76	0.84
1:A:749:GLN:N	1:A:753:THR:OG1	2.15	0.77
1:A:665:ASN:HD22	1:A:665:ASN:H	1.34	0.76
1:A:513:GLU:HG3	1:A:515:ILE:HG22	1.70	0.74
1:A:855:VAL:HG22	1:A:932:GLY:O	1.90	0.71
1:A:272:LYS:NZ	1:A:276:GLU:OE2	2.24	0.70
1:A:540:GLN:HG2	1:A:579:PRO:HA	1.74	0.70
1:A:364:GLU:HG2	1:A:367:SER:HB2	1.77	0.66
1:A:869:LEU:C	1:A:869:LEU:HD12	2.20	0.66
1:A:67:ALA:HB2	1:A:81:LYS:HD3	1.80	0.63
1:A:174:VAL:HG21	1:A:203:THR:HG21	1.81	0.61
1:A:117:SER:HB3	1:A:125:THR:HG21	1.83	0.60
1:A:149:ILE:HG22	1:A:501:LEU:HB3	1.83	0.60
1:A:665:ASN:HD22	1:A:665:ASN:N	1.99	0.58
1:A:194:ASN:ND2	1:A:483:LYS:HD2	2.19	0.58
1:A:194:ASN:HD22	1:A:483:LYS:HD2	1.68	0.58
1:A:561:SER:OG	1:A:564:SER:O	2.21	0.58
1:A:440:LYS:HB2	1:A:674:ASN:HB3	1.86	0.58
1:A:227:ASP:OD1	1:A:229:THR:HB	2.04	0.57
1:A:848:VAL:HG22	1:A:852:ARG:HD3	1.86	0.57
1:A:501:LEU:C	1:A:501:LEU:HD12	2.28	0.57
1:A:470:SER:HA	1:A:482:ILE:HD13	1.86	0.57
1:A:574:GLN:HE21	1:A:575:SER:H	1.51	0.57
1:A:246:ASN:N	1:A:246:ASN:OD1	2.35	0.57
1:A:859:VAL:HG11	1:A:890:LYS:HG3	1.87	0.56
1:A:156:LYS:HE2	6:A:1208:HOH:O	2.05	0.56
2:C:148:MET:HA	2:C:148:MET:HE2	1.89	0.55
1:A:555:ALA:HA	1:A:568:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:SER:OG	1:A:565:ASP:OD1	2.17	0.55
1:A:734:GLY:HA3	1:A:771:SER:O	2.08	0.53
1:A:626:SER:HA	1:A:651:LEU:O	2.08	0.53
1:A:281:LEU:HD11	1:A:296:TYR:CD1	2.43	0.53
1:A:290:VAL:HG12	1:A:294:LEU:HD22	1.90	0.52
1:A:317:LEU:HD23	1:A:317:LEU:N	2.23	0.52
1:A:341:ALA:HB1	1:A:378:PHE:HB3	1.92	0.52
1:A:231:GLN:O	1:A:268:TYR:OH	2.10	0.52
1:A:679:ILE:HD11	1:A:757:ASP:C	2.35	0.51
1:A:650:VAL:CG1	1:A:657:VAL:HG22	2.41	0.51
1:A:544:TYR:CD2	1:A:659:LEU:HD22	2.46	0.50
1:A:229:THR:HG22	1:A:230:LYS:HG3	1.93	0.50
1:A:390:SER:HA	1:A:394:GLN:O	2.11	0.50
1:A:650:VAL:HG12	1:A:657:VAL:HG22	1.94	0.50
1:A:723:ASP:HB3	1:A:725:LYS:H	1.77	0.50
1:A:554:SER:HB2	1:A:568:VAL:HG13	1.94	0.50
1:A:904:PHE:CG	1:A:905:LEU:N	2.80	0.50
1:A:457:PHE:CE1	3:B:3:SIA:H112	2.47	0.49
1:A:465:PHE:CB	1:A:515:ILE:HD13	2.41	0.49
1:A:173:GLU:O	1:A:177:GLY:HA2	2.12	0.49
1:A:454:ASN:OD1	1:A:467:VAL:HG23	2.12	0.49
1:A:682:SER:HG	1:A:761:SER:HB3	1.77	0.49
1:A:210:GLY:O	1:A:232:GLY:HA3	2.13	0.49
1:A:465:PHE:HB3	1:A:515:ILE:HD13	1.93	0.48
1:A:901:ARG:NE	6:A:1201:HOH:O	2.41	0.48
1:A:94:SER:HA	1:A:96:ASN:OD1	2.14	0.48
1:A:156:LYS:NZ	1:A:694:ASP:OD2	2.47	0.47
1:A:218:GLU:O	1:A:289:ASN:HA	2.15	0.47
1:A:451:GLY:HA3	1:A:485:ALA:O	2.15	0.47
1:A:558:SER:O	1:A:561:SER:HB3	2.15	0.47
1:A:158:THR:O	1:A:619:ALA:HB3	2.15	0.47
1:A:423:GLN:HG2	1:A:424:ALA:O	2.15	0.47
1:A:180:GLN:O	1:A:183:GLN:HG2	2.16	0.46
1:A:58:ALA:O	1:A:129:LYS:NZ	2.45	0.46
1:A:540:GLN:CG	1:A:579:PRO:HA	2.44	0.46
1:A:172:ARG:HH21	1:A:172:ARG:HG2	1.80	0.46
1:A:310:VAL:HA	1:A:323:TRP:O	2.16	0.46
1:A:453:ILE:O	1:A:467:VAL:HG22	2.16	0.46
1:A:669:LEU:HD23	1:A:679:ILE:HG23	1.98	0.46
1:A:230:LYS:HB3	1:A:231:GLN:HE22	1.81	0.45
2:C:144:ASN:HB3	2:C:148:MET:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:LYS:CB	1:A:231:GLN:NE2	2.80	0.45
1:A:896:LEU:HD12	1:A:896:LEU:C	2.42	0.45
1:A:159:LEU:HD23	1:A:619:ALA:HB1	1.98	0.45
1:A:439:TYR:CD2	1:A:678:PRO:HD3	2.52	0.44
1:A:558:SER:HA	1:A:566:THR:HB	2.00	0.44
1:A:132:LEU:HD22	1:A:323:TRP:CD1	2.53	0.44
1:A:99:PHE:CZ	1:A:108:PRO:HB3	2.53	0.43
1:A:357:GLU:OE2	1:A:379:THR:OG1	2.24	0.43
1:A:869:LEU:HD13	1:A:886:LEU:HB2	2.00	0.43
1:A:884:ALA:HA	1:A:901:ARG:O	2.19	0.43
2:C:180:TRP:HA	2:C:183:ILE:HG22	2.00	0.43
1:A:126:MET:HE1	1:A:129:LYS:HD2	2.00	0.43
1:A:310:VAL:HG21	1:A:395:ILE:HG13	2.00	0.42
1:A:677:ALA:O	1:A:678:PRO:C	2.62	0.42
1:A:649:VAL:HA	1:A:657:VAL:O	2.19	0.42
1:A:665:ASN:ND2	1:A:667:SER:OG	2.50	0.42
1:A:750:ASP:O	1:A:751:GLN:C	2.62	0.42
1:A:405:VAL:HG22	1:A:501:LEU:HD11	2.02	0.42
1:A:727:THR:O	1:A:728:PRO:C	2.62	0.42
1:A:364:GLU:HG2	1:A:367:SER:CB	2.46	0.42
1:A:138:SER:CB	1:A:565:ASP:OD1	2.68	0.42
1:A:658:ARG:NH1	1:A:746:GLU:OE2	2.47	0.42
1:A:679:ILE:HG12	1:A:759:ALA:HB2	2.01	0.42
1:A:866:PRO:HA	1:A:869:LEU:HG	2.02	0.42
1:A:317:LEU:N	1:A:317:LEU:CD2	2.82	0.42
1:A:44:ALA:HB3	1:A:799:ILE:HD13	2.01	0.42
1:A:210:GLY:H	1:A:232:GLY:HA3	1.85	0.42
1:A:792:LYS:HA	1:A:796:GLN:O	2.20	0.42
1:A:156:LYS:HE2	1:A:696:THR:OG1	2.20	0.41
1:A:391:VAL:CG2	1:A:631:ALA:HB1	2.50	0.41
1:A:554:SER:O	1:A:557:SER:HB3	2.20	0.41
1:A:668:LEU:HB3	1:A:680:SER:HB3	2.03	0.41
1:A:546:ILE:O	1:A:572:ASN:HB2	2.21	0.41
1:A:47:ILE:HG22	1:A:50:SER:HB3	2.03	0.41
1:A:120:ILE:HD13	1:A:918:GLY:O	2.20	0.41
1:A:398:SER:HA	1:A:544:TYR:HA	2.02	0.41
1:A:117:SER:HB3	1:A:125:THR:CG2	2.49	0.41
1:A:453:ILE:HG12	1:A:467:VAL:HG21	2.02	0.40
1:A:750:ASP:O	1:A:752:ALA:N	2.54	0.40
1:A:830:TYR:CE1	1:A:926:PRO:HD3	2.55	0.40
1:A:132:LEU:HD22	1:A:323:TRP:CG	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:ASP:CG	1:A:469:TYR:HA	2.46	0.40
1:A:77:PHE:CE2	1:A:113:GLU:HB2	2.56	0.40
1:A:574:GLN:HE21	1:A:575:SER:N	2.16	0.40
1:A:625:PHE:O	1:A:652:THR:HA	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	877/1053 (83%)	805 (92%)	62 (7%)	10 (1%)	11	46
2	C	1281/1444 (89%)	1176 (92%)	96 (8%)	9 (1%)	18	56
All	All	2158/2497 (86%)	1981 (92%)	158 (7%)	19 (1%)	16	51

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	SER
1	A	723	ASP
1	A	460	GLN
1	A	748	ASN
1	A	156	LYS
1	A	164	ALA
1	A	237	SER
1	A	529	SER
1	A	561	SER
2	C	643	SER
2	C	1092	THR
1	A	453	ILE
2	C	833	GLY
2	C	1346	PHE

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Mol	Chain	Res	Type
2	C	1187	LYS
2	C	641	ILE
2	C	822	PRO
2	C	593	THR
2	C	857	THR

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	775/914 (85%)	700 (90%)	75 (10%)	8	24
2	C	1122/1267 (89%)	1105 (98%)	17 (2%)	57	72
All	All	1897/2181 (87%)	1805 (95%)	92 (5%)	24	44

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	52	SER
1	A	69	GLN
1	A	79	LEU
1	A	81	LYS
1	A	90	GLU
1	A	104	GLN
1	A	106	GLU
1	A	125	THR
1	A	149	ILE
1	A	163	LEU
1	A	169	LYS
1	A	172	ARG
1	A	180	GLN
1	A	182	ASN
1	A	186	SER
1	A	190	GLN
1	A	229	THR

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Mol	Chain	Res	Type
1	A	231	GLN
1	A	237	SER
1	A	241	SER
1	A	246	ASN
1	A	249	LYS
1	A	264	THR
1	A	275	LEU
1	A	277	SER
1	A	286	THR
1	A	294	LEU
1	A	302	LEU
1	A	317	LEU
1	A	364	GLU
1	A	368	THR
1	A	376	LYS
1	A	377	SER
1	A	390	SER
1	A	393	ASP
1	A	409	SER
1	A	421	LYS
1	A	437	THR
1	A	441	ASP
1	A	459	ILE
1	A	466	VAL
1	A	469	TYR
1	A	470	SER
1	A	492	GLN
1	A	493	LYS
1	A	495	THR
1	A	500	SER
1	A	521	LEU
1	A	528	LYS
1	A	530	ASN
1	A	533	ARG
1	A	536	SER
1	A	537	ARG
1	A	540	GLN
1	A	543	VAL
1	A	561	SER
1	A	574	GLN
1	A	576	ARG
1	A	643	LYS

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Mol	Chain	Res	Type
1	A	657	VAL
1	A	665	ASN
1	A	676	ILE
1	A	679	ILE
1	A	680	SER
1	A	684	THR
1	A	725	LYS
1	A	765	VAL
1	A	766	GLN
1	A	792	LYS
1	A	793	ASP
1	A	855	VAL
1	A	863	ILE
1	A	869	LEU
1	A	894	THR
2	C	243	ASP
2	C	251	LYS
2	C	401	ASN
2	C	503	TRP
2	C	622	PHE
2	C	638	GLU
2	C	650	LEU
2	C	725	ARG
2	C	752	THR
2	C	778	SER
2	C	819	VAL
2	C	876	LYS
2	C	887	ILE
2	C	890	THR
2	C	1030	THR
2	C	1231	THR
2	C	1344	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	182	ASN
1	A	183	GLN
1	A	187	GLN
1	A	189	ASN
1	A	194	ASN

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Mol	Chain	Res	Type
1	A	200	ASN
1	A	231	GLN
1	A	250	ASN
1	A	285	GLN
1	A	293	GLN
1	A	330	GLN
1	A	352	GLN
1	A	369	ASN
1	A	492	GLN
1	A	499	ASN
1	A	574	GLN
1	A	665	ASN
1	A	674	ASN
1	A	770	GLN
2	C	199	ASN
2	C	215	ASN
2	C	462	ASN
2	C	631	ASN
2	C	633	ASN
2	C	634	ASN
2	C	665	GLN
2	C	675	ASN
2	C	716	ASN
2	C	727	ASN
2	C	735	ASN
2	C	745	ASN
2	C	779	ASN
2	C	815	ASN
2	C	837	GLN
2	C	840	ASN
2	C	849	GLN
2	C	862	ASN
2	C	908	ASN
2	C	1003	ASN
2	C	1235	ASN
2	C	1240	GLN

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 ⓘ

3 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	ASO	B	1	3	11,11,11	1.51	3 (27%)	15,15,15	0.83	0
3	GAL	B	2	3	11,11,12	1.34	2 (18%)	15,15,17	0.79	0
3	SIA	B	3	3	20,20,21	2.02	2 (10%)	21,28,31	1.34	3 (14%)

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	ASO	B	1	3	-	0/2/19/19	0/1/1/1
3	GAL	B	2	3	-	0/2/19/22	0/1/1/1
3	SIA	B	3	3	-	0/18/34/38	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3	SIA	C2-C1	7.45	1.61	1.52
3	B	3	SIA	C4-C5	2.80	1.55	1.53
3	B	1	ASO	O5-C1	2.73	1.48	1.43
3	B	2	GAL	O5-C5	2.68	1.48	1.43
3	B	1	ASO	O5-C5	2.59	1.48	1.43
3	B	1	ASO	O4-C4	2.47	1.49	1.43
3	B	2	GAL	O5-C1	2.03	1.47	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3	SIA	O6-C2-C1	3.26	113.88	107.72
3	B	3	SIA	O6-C2-C3	-2.61	107.05	110.56
3	B	3	SIA	O1B-C1-C2	2.31	118.71	112.71

There are no chirality outliers.

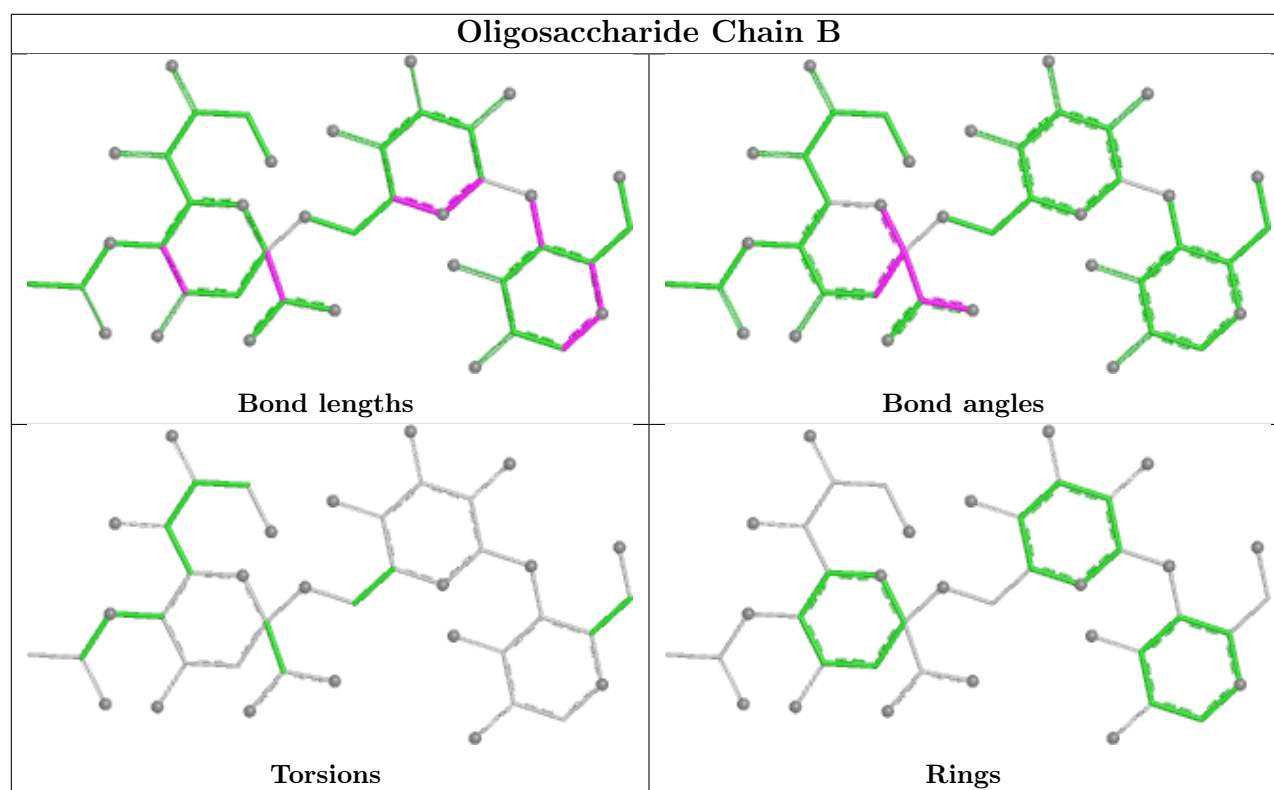
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	3	SIA	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](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	PO4	A	1102	-	4,4,4	0.73	0	6,6,6	0.59	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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.

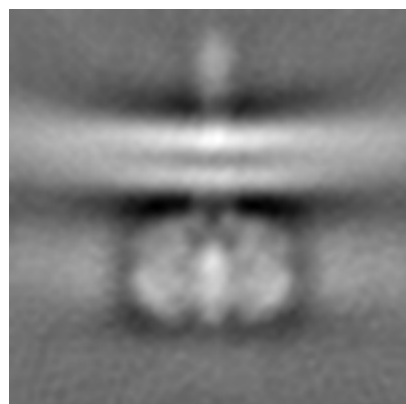
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17592. These allow visual inspection of the internal detail of the map and identification of artifacts.

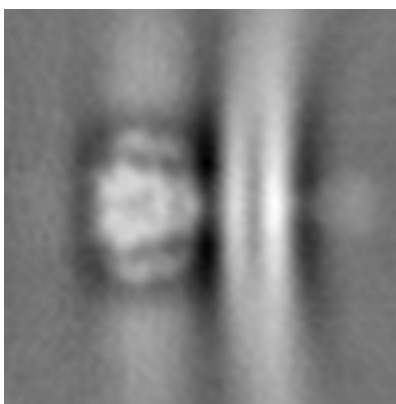
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

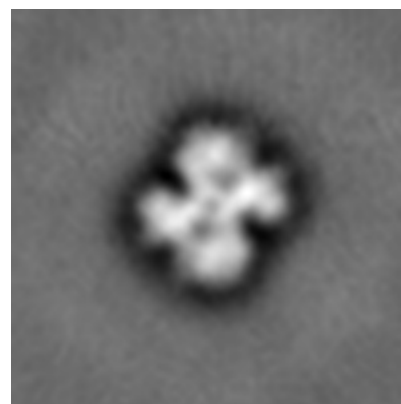
6.1.1 Primary map



X

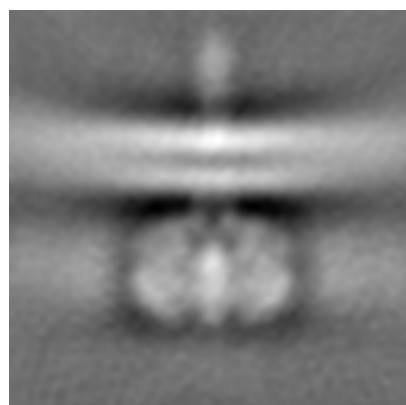


Y

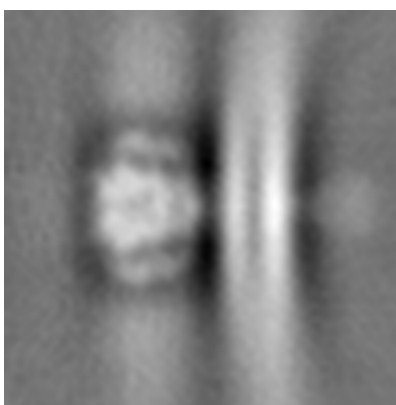


Z

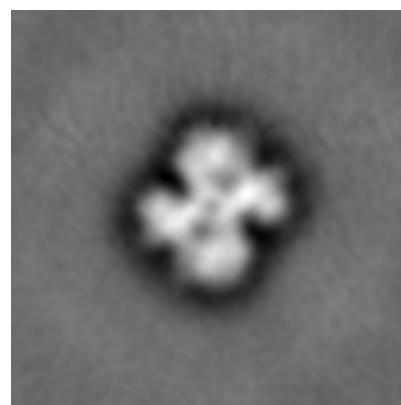
6.1.2 Raw map



X



Y

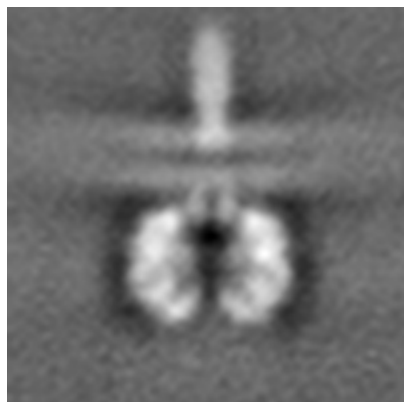


Z

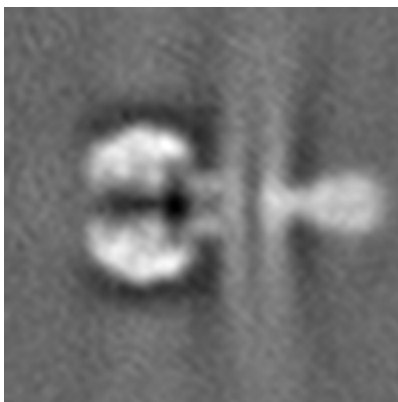
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

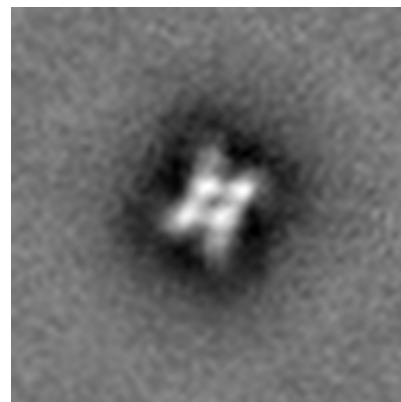
6.2.1 Primary map



X Index: 32

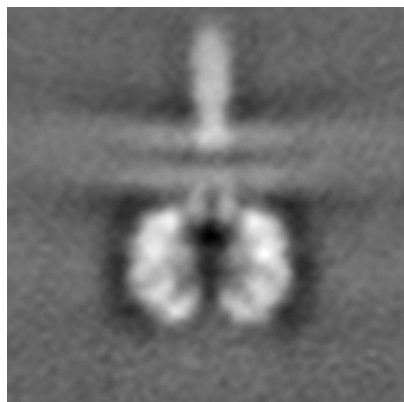


Y Index: 32

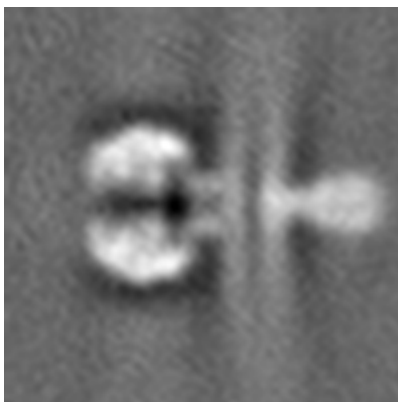


Z Index: 32

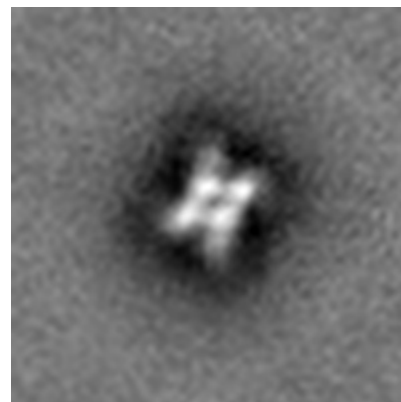
6.2.2 Raw map



X Index: 32



Y Index: 32

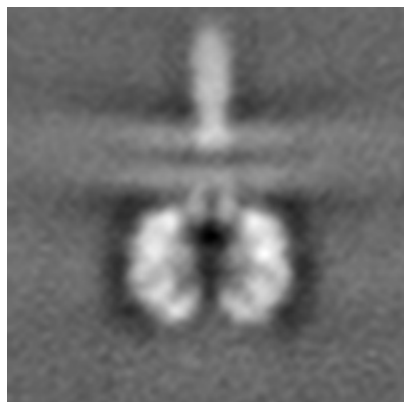


Z Index: 32

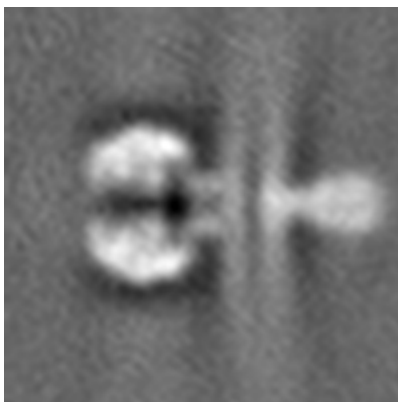
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

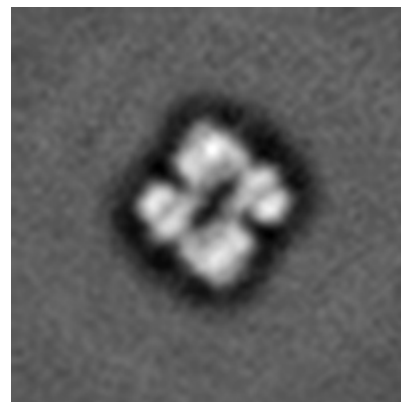
6.3.1 Primary map



X Index: 32

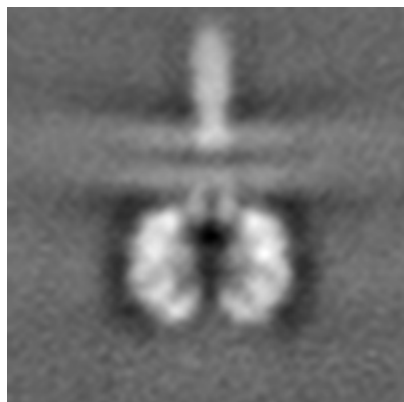


Y Index: 32

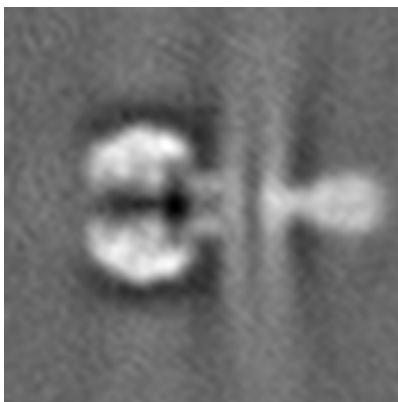


Z Index: 24

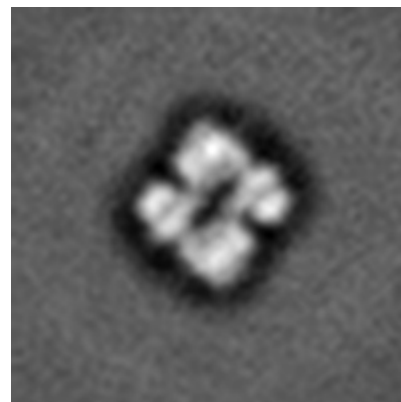
6.3.2 Raw map



X Index: 32



Y Index: 32

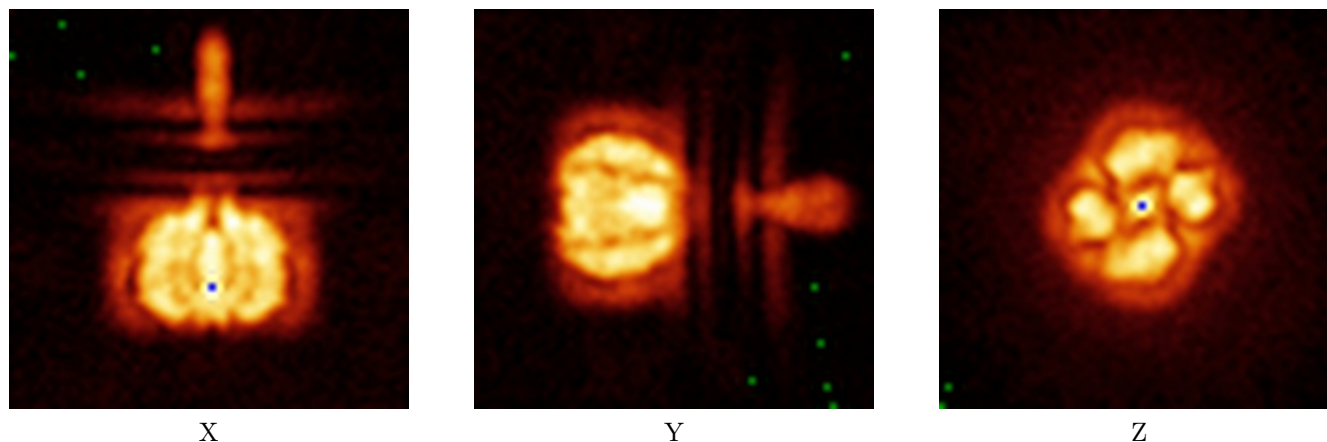


Z Index: 24

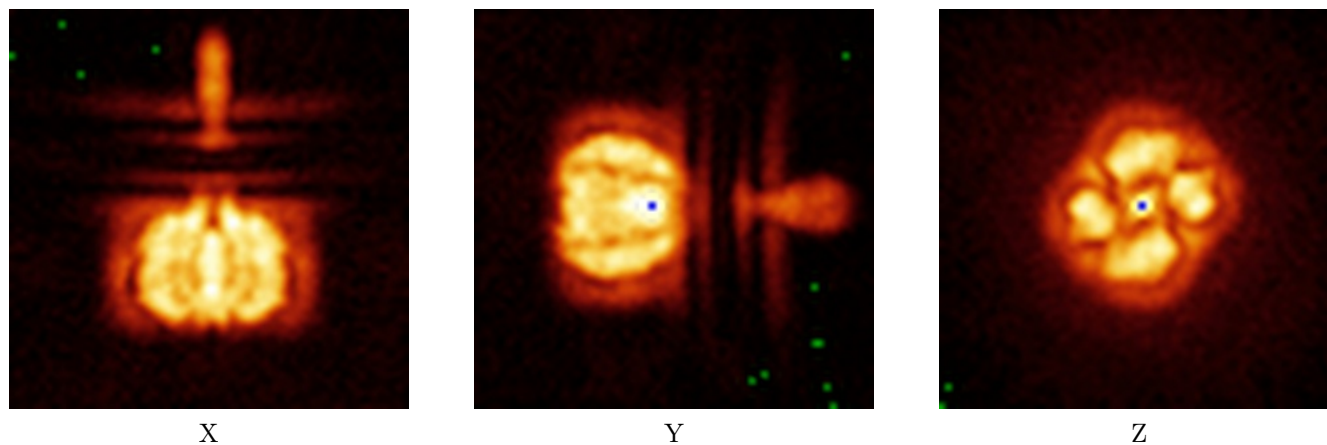
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



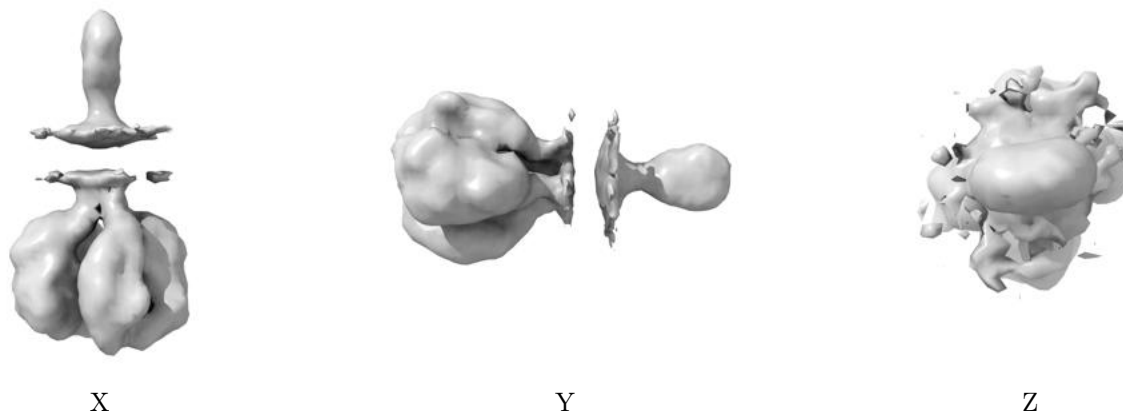
6.4.2 Raw map



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.

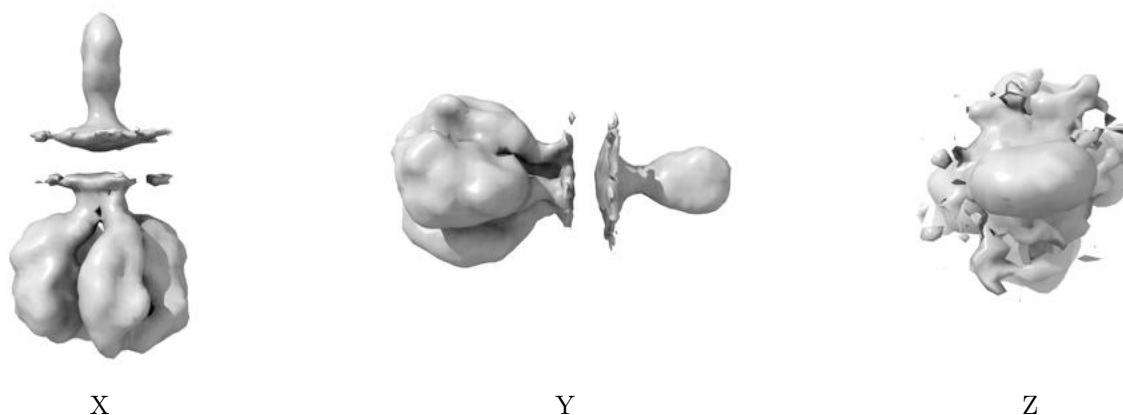
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.000229. 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.

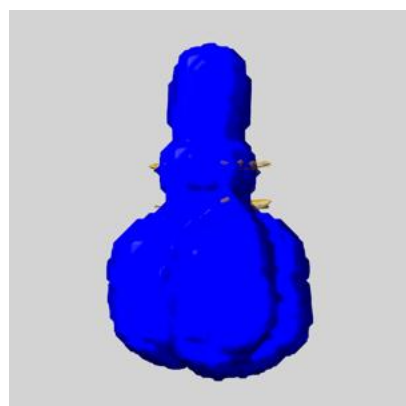
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

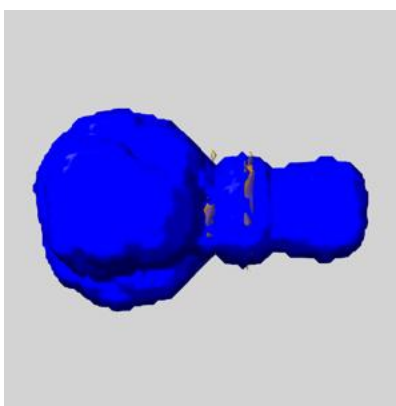
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

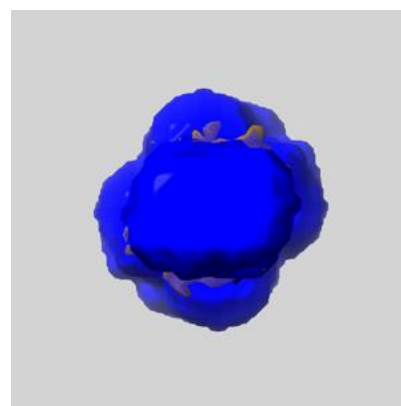
6.6.1 emd_17592_msk_1.map [i](#)



X



Y

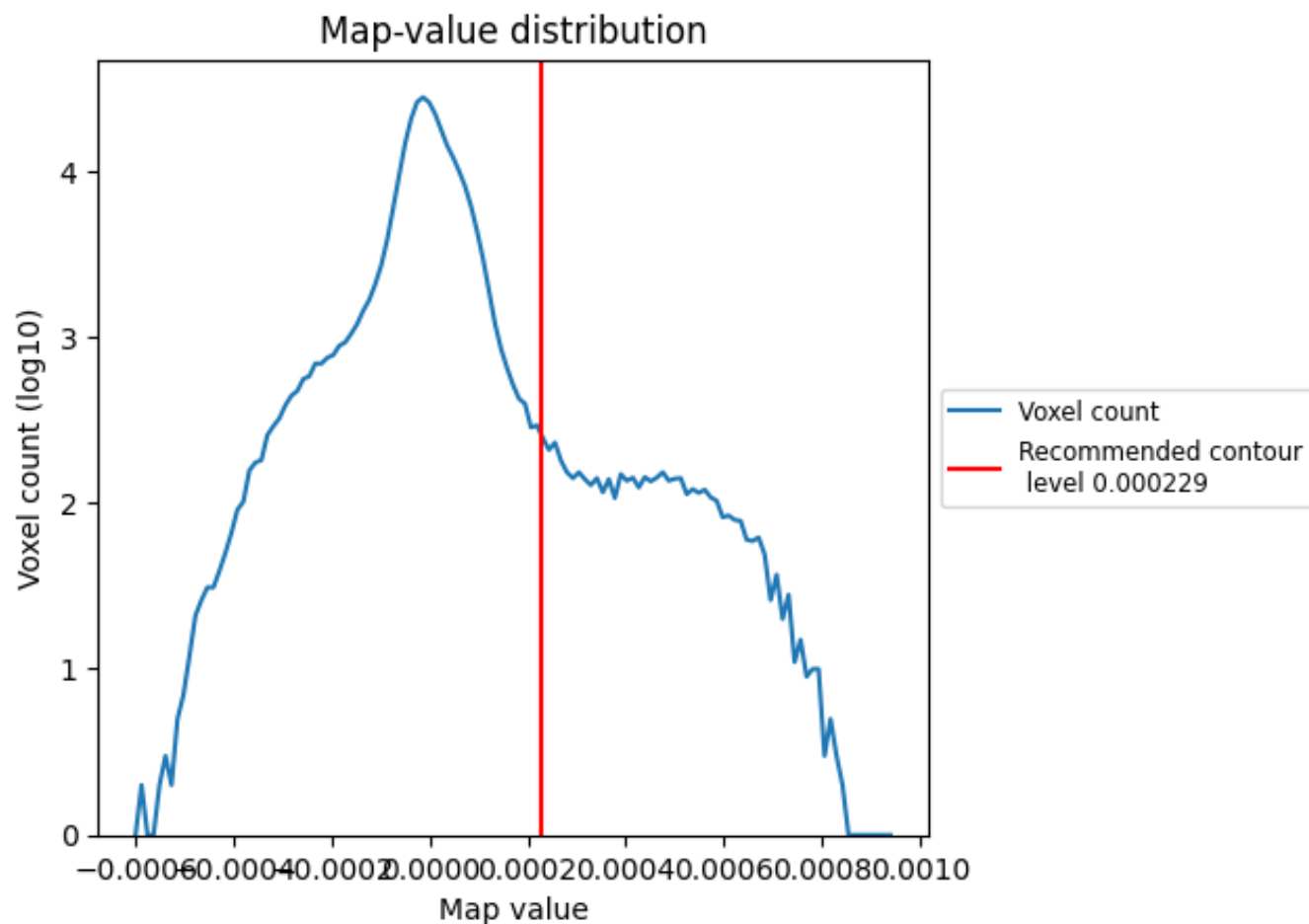


Z

7 Map analysis [i](#)

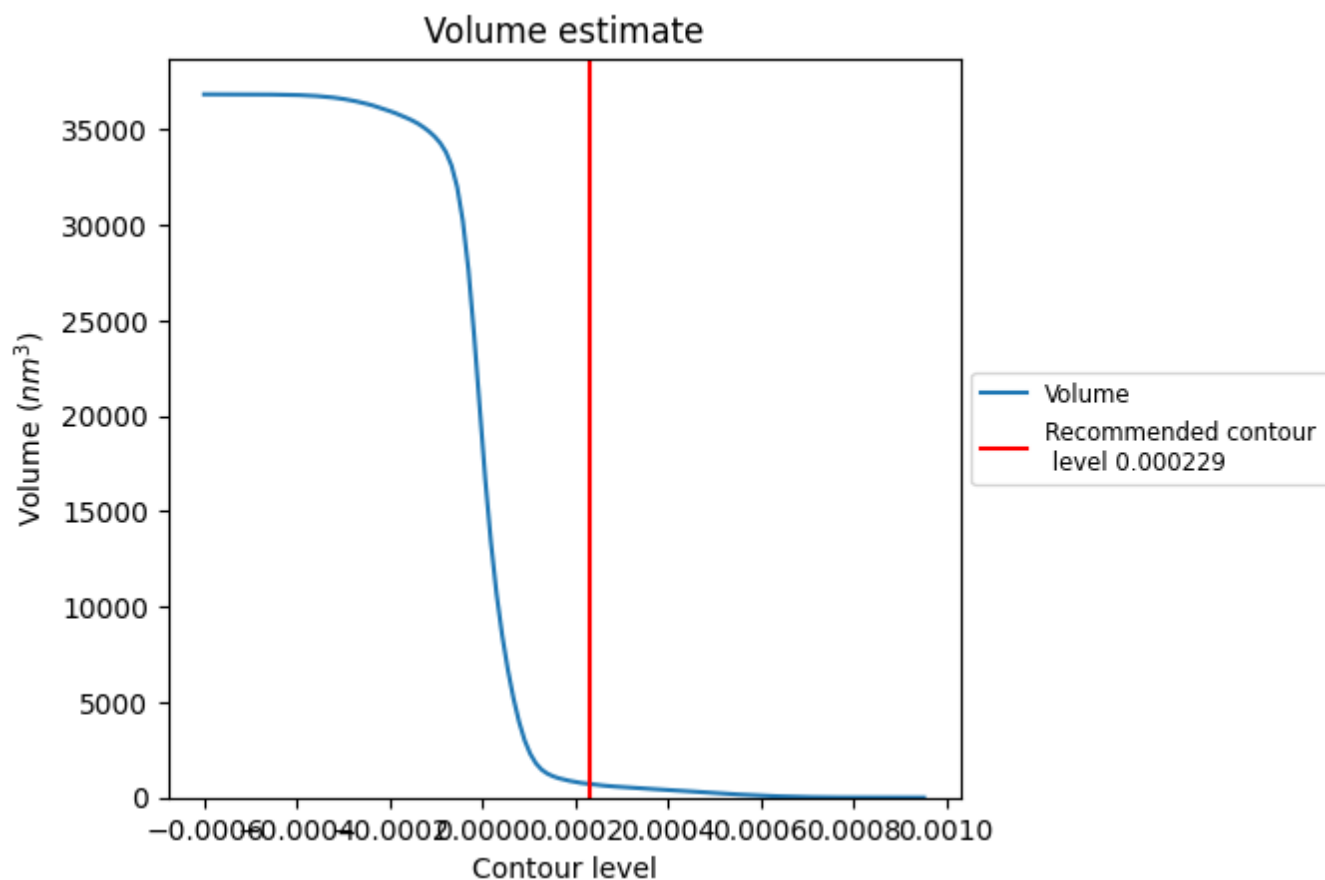
This section contains the results of statistical analysis of the map.

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.

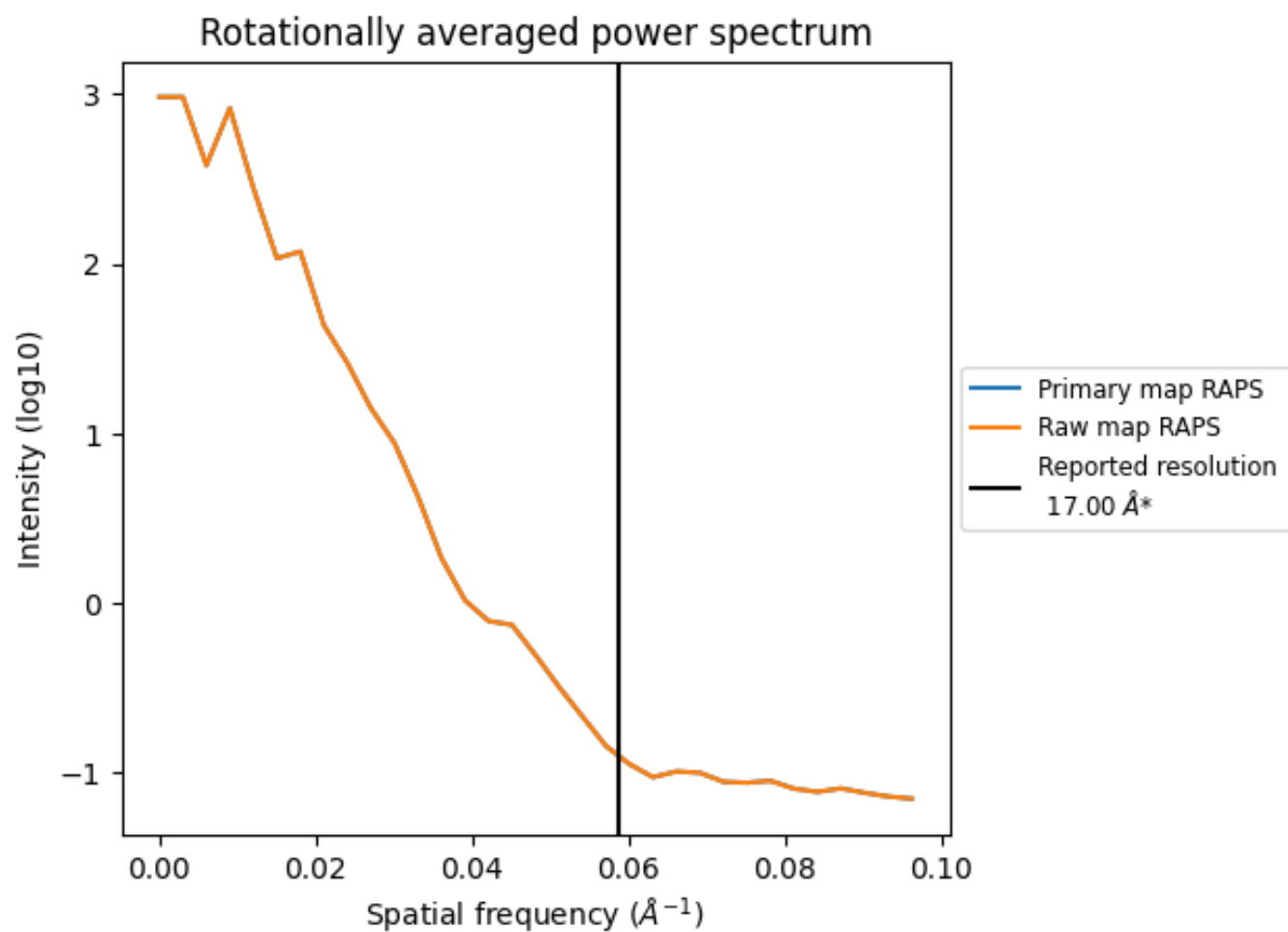
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 717 nm³; this corresponds to an approximate mass of 648 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 ⓘ

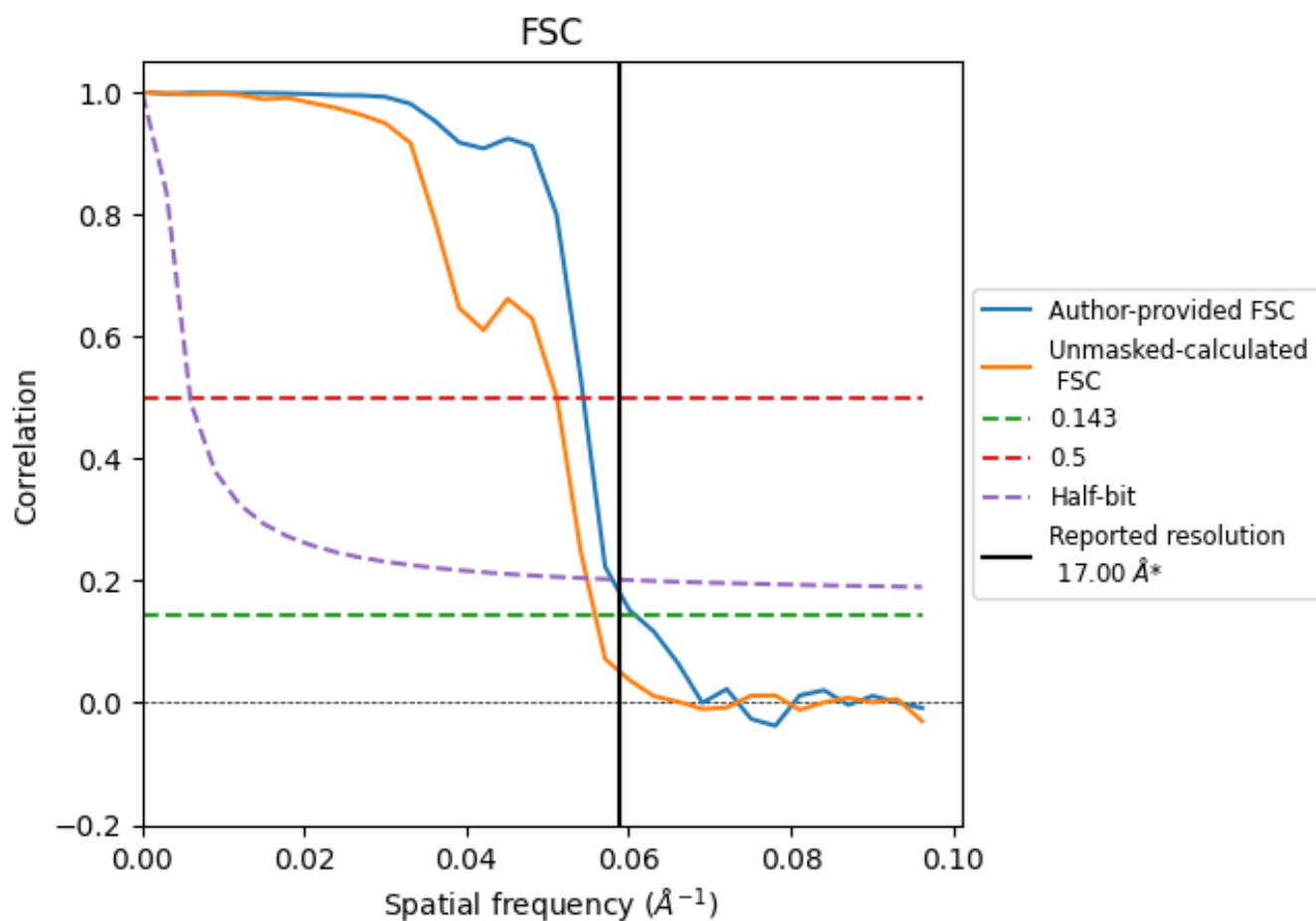


*Reported resolution corresponds to spatial frequency of 0.059 Å⁻¹

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 0.059 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

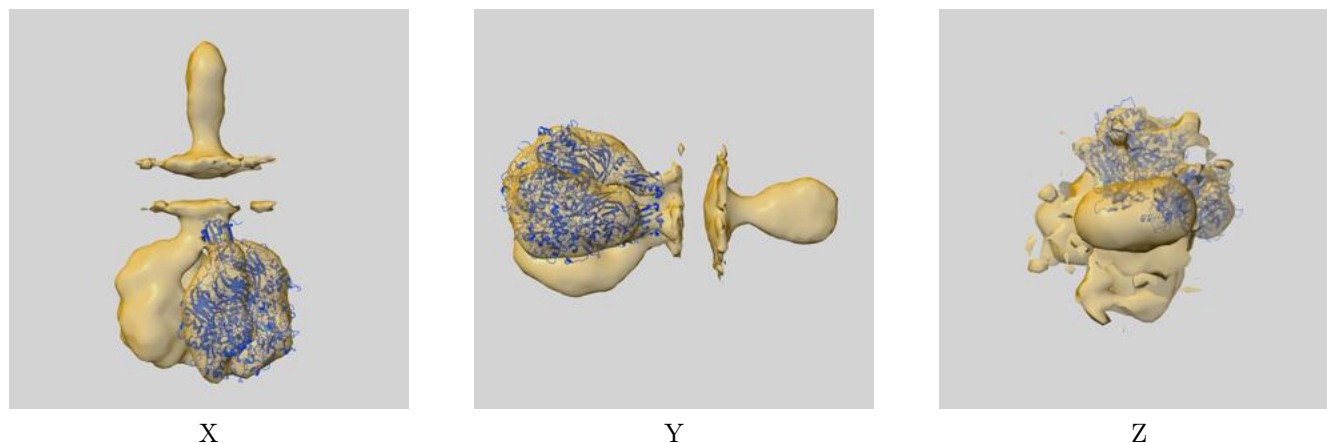
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	17.00	-	-
Author-provided FSC curve	16.42	18.38	17.24
Unmasked-calculated*	17.89	19.57	18.25

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

9 Map-model fit [i](#)

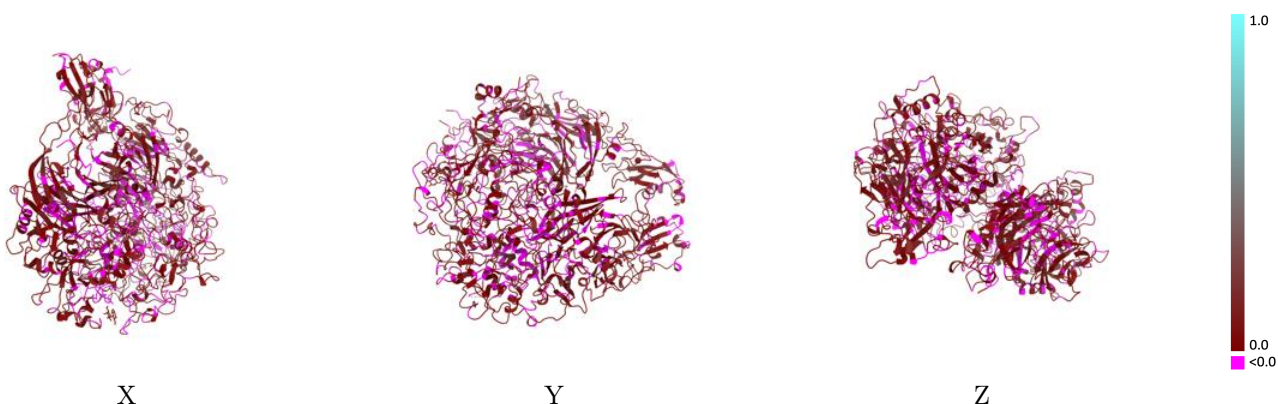
This section contains information regarding the fit between EMDB map EMD-17592 and PDB model 8PC0. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



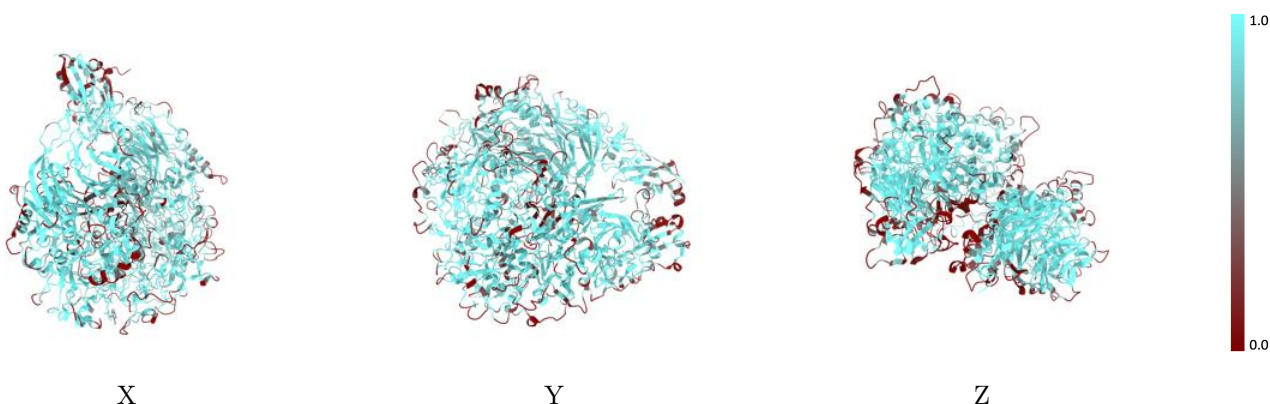
The images above show the 3D surface view of the map at the recommended contour level 0.000229 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



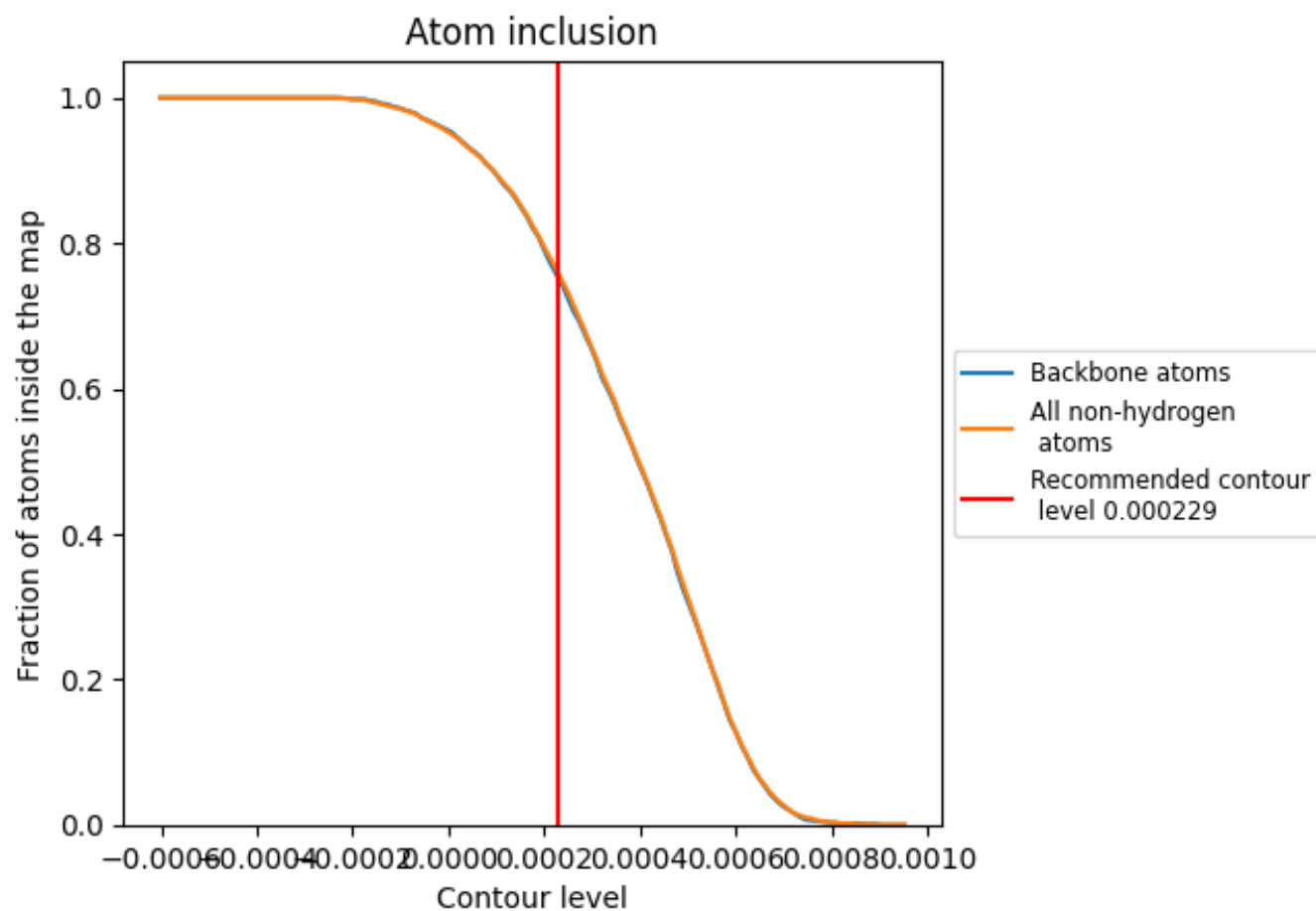
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.000229).

9.4 Atom inclusion ⓘ



At the recommended contour level, 75% of all backbone atoms, 76% 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 (0.000229) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7600	0.0480
A	0.7860	0.0530
B	0.8330	-0.0070
C	0.7440	0.0450

