



Full wwPDB EM Validation Report ⓘ

Jul 14, 2026 – 05:27 pm BST

PDB ID : 28JL / pdb_000028jl
EMDB ID : EMD-56543
Title : Lid-engaged Complex: BAM bound BepA with lid engaged in membrane
Authors : Fenn, K.L.; Ranson, N.A.
Deposited on : 2026-02-03
Resolution : 3.60 Å (reported)
Based on initial models : ., 5LJO

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:

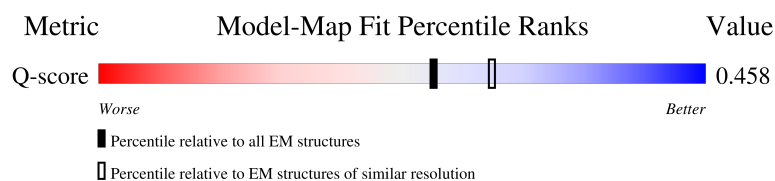
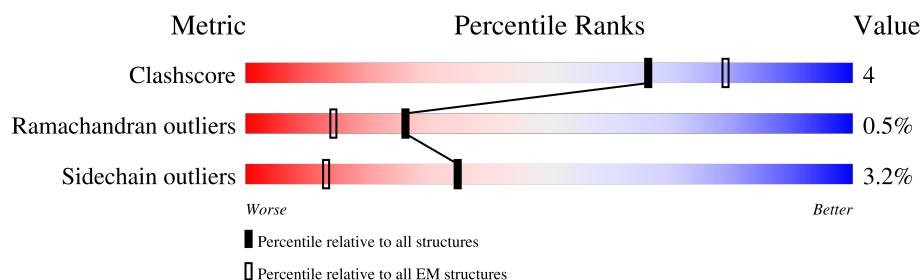
EMDB validation analysis : 0.0.1.dev133
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

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 3.60 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	12797 (3.10 - 4.10)

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	790	
2	B	373	
3	C	320	
4	D	226	

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Mol	Chain	Length	Quality of chain
5	E	104	5% 81% 8% 12%
6	F	492	89% 7%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17104 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 Outer membrane protein assembly factor BamA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	774	Total	C	N	O	S	0	0
			6050	3819	1028	1187	16		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	723	CYS	PRO	engineered mutation	UNP P0A940

- Molecule 2 is a protein called Outer membrane protein assembly factor BamB.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	365	Total	C	N	O	S	0	0
			2737	1713	469	549	6		

- Molecule 3 is a protein called Outer membrane protein assembly factor BamC.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	319	Total	C	N	O	S	0	0
			2274	1407	409	451	7		

- Molecule 4 is a protein called Outer membrane protein assembly factor BamD.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	217	Total	C	N	O	S	0	0
			1755	1106	308	334	7		

- Molecule 5 is a protein called Outer membrane protein assembly factor BamE.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	92	Total	C	N	O	S	0	0
			718	451	124	140	3		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	114	GLY	-	expression tag	UNP P0A937
E	115	GLY	-	expression tag	UNP P0A937
E	116	HIS	-	expression tag	UNP P0A937
E	117	HIS	-	expression tag	UNP P0A937
E	118	HIS	-	expression tag	UNP P0A937
E	119	HIS	-	expression tag	UNP P0A937
E	120	HIS	-	expression tag	UNP P0A937
E	121	HIS	-	expression tag	UNP P0A937
E	122	HIS	-	expression tag	UNP P0A937
E	123	HIS	-	expression tag	UNP P0A937

- Molecule 6 is a protein called Beta-barrel assembly-enhancing protease.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	459	Total	C	N	O	S	0	0
			3569	2209	657	687	16		

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	186	CYS	LEU	engineered mutation	UNP P66948
F	488	GLY	-	expression tag	UNP P66948
F	489	SER	-	expression tag	UNP P66948
F	490	SER	-	expression tag	UNP P66948
F	491	ALA	-	expression tag	UNP P66948
F	492	TRP	-	expression tag	UNP P66948
F	493	SER	-	expression tag	UNP P66948
F	494	HIS	-	expression tag	UNP P66948
F	495	PRO	-	expression tag	UNP P66948
F	496	GLN	-	expression tag	UNP P66948
F	497	PHE	-	expression tag	UNP P66948
F	498	GLU	-	expression tag	UNP P66948
F	499	LYS	-	expression tag	UNP P66948
F	500	GLY	-	expression tag	UNP P66948
F	501	GLY	-	expression tag	UNP P66948
F	502	GLY	-	expression tag	UNP P66948
F	503	SER	-	expression tag	UNP P66948
F	504	GLY	-	expression tag	UNP P66948
F	505	GLY	-	expression tag	UNP P66948
F	506	GLY	-	expression tag	UNP P66948
F	507	SER	-	expression tag	UNP P66948
F	508	GLY	-	expression tag	UNP P66948
F	509	GLY	-	expression tag	UNP P66948

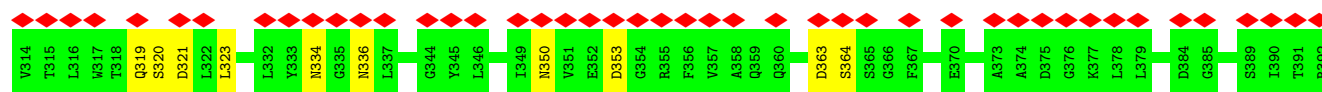
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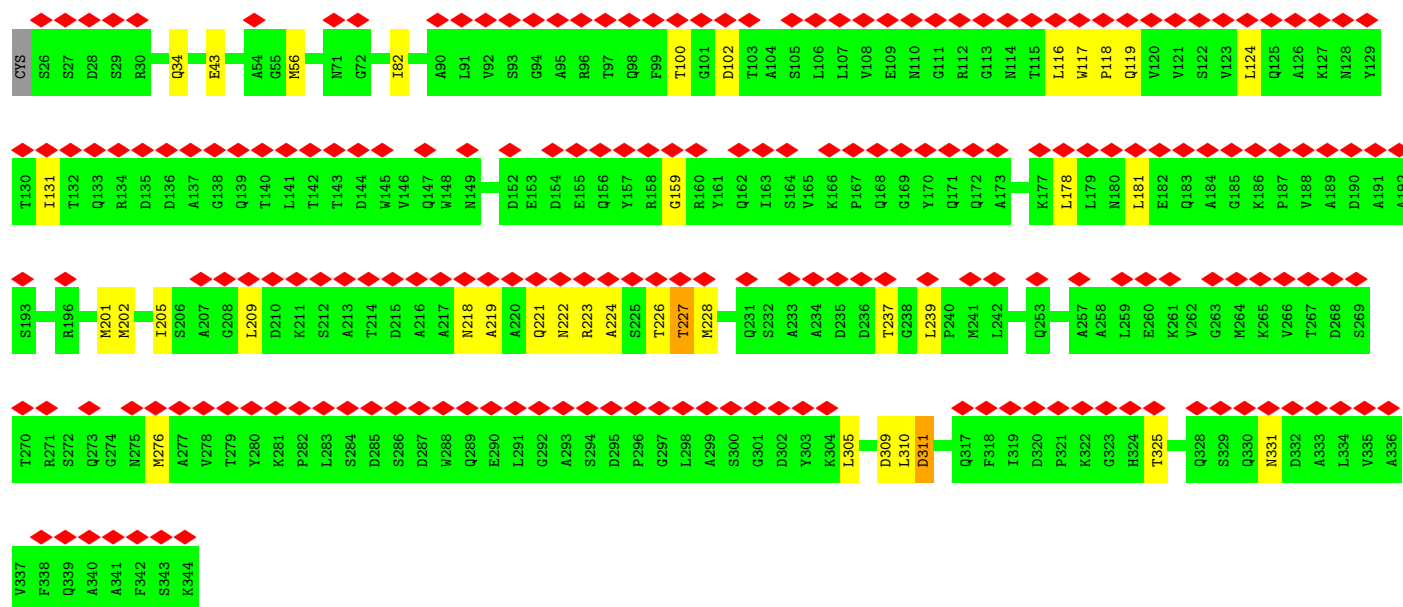
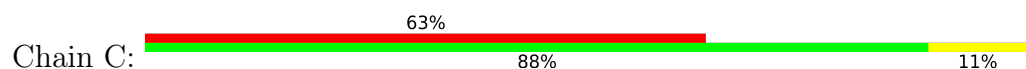
Chain	Residue	Modelled	Actual	Comment	Reference
F	510	SER	-	expression tag	UNP P66948
F	511	ALA	-	expression tag	UNP P66948
F	512	TRP	-	expression tag	UNP P66948
F	513	SER	-	expression tag	UNP P66948
F	514	HIS	-	expression tag	UNP P66948
F	515	PRO	-	expression tag	UNP P66948
F	516	GLN	-	expression tag	UNP P66948
F	517	PHE	-	expression tag	UNP P66948
F	518	GLU	-	expression tag	UNP P66948
F	519	LYS	-	expression tag	UNP P66948

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
7	F	1	Total 1	Zn 1	0



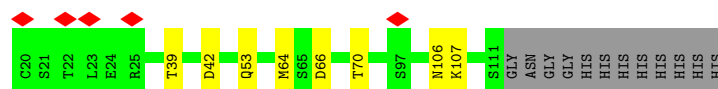
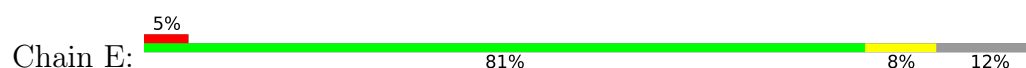
- Molecule 3: Outer membrane protein assembly factor BamC



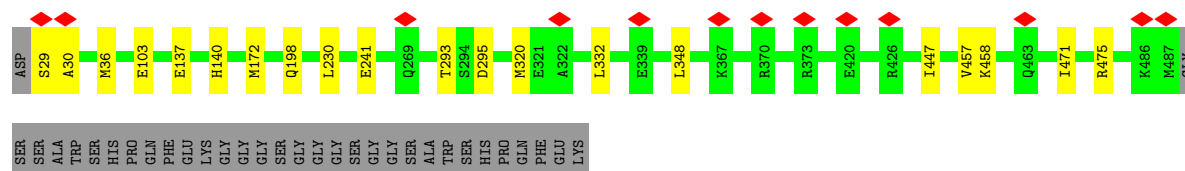
- Molecule 4: Outer membrane protein assembly factor BamD



- Molecule 5: Outer membrane protein assembly factor BamE



- Molecule 6: Beta-barrel assembly-enhancing protease



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	126000	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	900	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.881	Depositor
Minimum map value	-0.542	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.145	Depositor
Map size (Å)	228.0, 228.0, 228.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.76, 0.76, 0.76	Depositor

5 Model quality

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.35	0/6183	0.51	3/8388 (0.0%)
2	B	0.22	0/2784	0.30	0/3796
3	C	0.43	0/2311	0.61	0/3156
4	D	0.23	0/1795	0.38	0/2439
5	E	0.11	0/733	0.22	0/1000
6	F	0.16	0/3629	0.35	0/4912
All	All	0.29	0/17435	0.44	3/23691 (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	A	0	5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	763	PRO	N-CA-C	-5.57	106.44	114.18
1	A	678	GLN	N-CA-C	-5.55	106.51	113.72
1	A	265	GLN	CA-C-O	-5.02	116.23	121.55

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	321	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	353	ARG	Sidechain
1	A	632	ARG	Sidechain
1	A	634	ARG	Sidechain
1	A	767	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	A	6050	0	5747	55	0
2	B	2737	0	2673	19	0
3	C	2274	0	2119	32	0
4	D	1755	0	1694	9	0
5	E	718	0	699	5	0
6	F	3569	0	3502	14	0
7	F	1	0	0	0	0
All	All	17104	0	16434	129	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 4.

All (129) 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)
3:C:159:GLY:HA3	3:C:181:LEU:HD23	1.56	0.86
6:F:172:MET:HA	6:F:172:MET:HE3	1.58	0.86
1:A:114:LEU:HD21	1:A:137:LEU:HD21	1.65	0.76
3:C:178:LEU:HD11	3:C:181:LEU:HG	1.67	0.75
6:F:103:GLU:OE2	6:F:241:GLU:HG3	1.87	0.74
5:E:66:ASP:OD2	5:E:70:THR:OG1	2.09	0.71
3:C:311:ASP:OD1	6:F:475:ARG:NH2	2.24	0.70
3:C:178:LEU:HD21	3:C:181:LEU:CD1	2.22	0.70
1:A:369:MET:HA	1:A:369:MET:HE2	1.73	0.69
5:E:53:GLN:N	5:E:53:GLN:OE1	2.26	0.69
4:D:26:GLU:OE1	4:D:26:GLU:N	2.26	0.68
1:A:393:GLY:O	1:A:421:ARG:NH1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:GLU:OE2	1:A:391:ARG:NH1	2.27	0.67
1:A:377:LEU:HD13	1:A:378:GLY:N	2.10	0.66
3:C:43:GLU:N	3:C:43:GLU:OE1	2.27	0.66
3:C:178:LEU:HD11	3:C:181:LEU:CG	2.26	0.66
3:C:159:GLY:CA	3:C:181:LEU:HD23	2.25	0.66
1:A:160:ARG:NH1	3:C:237:THR:O	2.30	0.65
1:A:77:VAL:HG13	1:A:84:LEU:HD21	1.79	0.64
1:A:746:ASP:OD1	1:A:747:THR:N	2.31	0.63
2:B:300:ASP:OD1	2:B:304:ARG:N	2.31	0.63
3:C:159:GLY:HA3	3:C:181:LEU:CD2	2.28	0.63
1:A:470:GLU:OE2	1:A:488:ARG:NH1	2.32	0.62
2:B:79:GLY:HA2	2:B:111:LEU:HD11	1.82	0.62
3:C:116:LEU:HD11	3:C:209:LEU:HD21	1.81	0.62
3:C:178:LEU:HD21	3:C:181:LEU:HD11	1.81	0.62
3:C:226:THR:O	3:C:227:THR:C	2.43	0.62
1:A:171:GLU:HB2	2:B:27:GLU:HG3	1.81	0.61
2:B:88:ASP:N	2:B:88:ASP:OD1	2.34	0.60
1:A:211:ASP:N	1:A:211:ASP:OD1	2.35	0.59
4:D:212:ARG:NH2	4:D:241:ASN:O	2.35	0.59
1:A:684:PRO:O	1:A:685:ASP:C	2.45	0.59
3:C:239:LEU:HD13	3:C:325:THR:HG21	1.85	0.59
4:D:92:ILE:HG23	4:D:108:VAL:HG13	1.86	0.58
1:A:399:ASP:N	1:A:399:ASP:OD1	2.36	0.58
1:A:264:ASP:N	1:A:264:ASP:OD1	2.38	0.57
1:A:623:ASP:OD1	1:A:624:ASP:N	2.37	0.57
3:C:276:MET:N	3:C:305:LEU:O	2.37	0.57
1:A:332:ASP:OD1	1:A:332:ASP:N	2.36	0.56
1:A:615:THR:HG23	1:A:633:THR:HB	1.87	0.56
1:A:250:ASP:OD1	1:A:250:ASP:N	2.38	0.56
1:A:655:GLY:O	1:A:659:THR:OG1	2.24	0.56
5:E:39:THR:OG1	5:E:42:ASP:OD2	2.22	0.55
3:C:56:MET:HE3	3:C:56:MET:HA	1.87	0.55
1:A:672:ALA:O	1:A:702:SER:OG	2.20	0.55
3:C:178:LEU:HD21	3:C:181:LEU:HD12	1.89	0.55
1:A:663:PHE:HD2	1:A:668:ILE:HD11	1.72	0.54
2:B:350:ASN:ND2	2:B:353:ASP:OD1	2.41	0.54
1:A:365:LEU:HD21	1:A:395:PHE:CZ	2.43	0.54
3:C:119:GLN:HB3	3:C:209:LEU:HD11	1.91	0.52
2:B:334:ASN:O	2:B:336:ASN:ND2	2.42	0.52
2:B:80:LEU:HD21	2:B:82:LYS:HE3	1.91	0.52
3:C:178:LEU:CD2	3:C:181:LEU:HD11	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:595:LEU:C	1:A:595:LEU:HD23	2.35	0.52
3:C:124:LEU:HD12	3:C:124:LEU:O	2.10	0.52
3:C:202:MET:O	3:C:205:ILE:HG22	2.09	0.52
4:D:117:MET:CE	4:D:170:LEU:HD21	2.41	0.51
4:D:92:ILE:HG21	4:D:112:ARG:HB2	1.93	0.51
5:E:106:ASN:OD1	5:E:107:LYS:N	2.44	0.51
2:B:319:GLN:OE1	2:B:320:SER:N	2.44	0.51
4:D:125:GLN:OE1	4:D:132:ARG:NH1	2.39	0.50
2:B:363:ASP:OD1	2:B:364:SER:N	2.45	0.50
2:B:202:THR:HG22	2:B:221:MET:HE1	1.94	0.49
1:A:197:PHE:HZ	1:A:226:LEU:HD12	1.77	0.49
2:B:192:LEU:HD22	6:F:36:MET:HE2	1.95	0.49
1:A:660:VAL:O	1:A:661:ARG:C	2.56	0.48
1:A:752:SER:OG	1:A:753:GLN:NE2	2.42	0.48
1:A:568:ASP:OD1	1:A:568:ASP:N	2.47	0.47
3:C:34:GLN:O	3:C:82:ILE:HD11	2.15	0.47
6:F:293:THR:OG1	6:F:295:ASP:OD1	2.32	0.47
2:B:319:GLN:OE1	2:B:321:ASP:N	2.43	0.47
1:A:239:ASN:OD1	1:A:240:ILE:N	2.48	0.47
1:A:505:THR:HG1	1:A:537:SER:HG	1.55	0.47
3:C:117:TRP:HB3	3:C:118:PRO:HD3	1.96	0.47
6:F:332:LEU:CD1	6:F:348:LEU:HD23	2.45	0.47
1:A:560:ASP:HA	1:A:679:ALA:HB2	1.96	0.47
1:A:328:ILE:HA	1:A:335:VAL:HG12	1.97	0.47
4:D:195:VAL:HG11	5:E:64:MET:HE3	1.97	0.46
3:C:309:ASP:C	3:C:310:LEU:HD22	2.41	0.46
2:B:111:LEU:HD12	2:B:111:LEU:H	1.80	0.46
3:C:116:LEU:CD1	3:C:209:LEU:HD21	2.44	0.46
1:A:104:SER:OG	1:A:171:GLU:OE1	2.33	0.46
1:A:232:ASP:O	1:A:297:ASN:ND2	2.49	0.46
2:B:272:LEU:C	2:B:272:LEU:HD23	2.41	0.45
6:F:103:GLU:OE2	6:F:241:GLU:CG	2.60	0.45
3:C:124:LEU:HD21	3:C:131:ILE:HD11	1.99	0.45
1:A:85:LEU:HD23	1:A:86:VAL:N	2.32	0.45
1:A:605:ASP:OD1	1:A:605:ASP:N	2.50	0.45
2:B:271:ASP:OD2	2:B:274:SER:N	2.51	0.44
6:F:320:MET:C	6:F:320:MET:SD	3.00	0.44
6:F:172:MET:HA	6:F:172:MET:CE	2.40	0.44
1:A:405:VAL:HG11	1:A:411:GLN:HB2	2.00	0.44
1:A:505:THR:OG1	1:A:537:SER:OG	2.26	0.44
3:C:227:THR:O	3:C:228:MET:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:VAL:HG11	1:A:252:LYS:HZ2	1.83	0.43
4:D:187:GLU:OE2	4:D:187:GLU:O	2.35	0.43
1:A:268:LEU:HD12	1:A:335:VAL:HG23	2.00	0.43
1:A:185:ASN:ND2	1:A:188:PHE:O	2.51	0.43
1:A:594:ASN:O	1:A:613:LEU:HD12	2.18	0.43
3:C:102:ASP:N	3:C:102:ASP:OD1	2.51	0.43
1:A:805:ASN:OD1	1:A:806:ILE:N	2.51	0.43
4:D:148:SER:OG	4:D:152:ARG:NH2	2.52	0.43
6:F:137:GLU:OE2	6:F:137:GLU:HA	2.18	0.43
2:B:281:ARG:NH1	2:B:313:GLY:O	2.51	0.42
1:A:351:LYS:HB2	1:A:351:LYS:HE2	1.31	0.42
3:C:201:MET:SD	3:C:201:MET:O	2.78	0.42
3:C:218:ASN:CG	3:C:222:ASN:H	2.28	0.42
3:C:223:ARG:O	3:C:224:ALA:C	2.63	0.42
3:C:226:THR:O	3:C:228:MET:N	2.52	0.42
6:F:457:VAL:HG12	6:F:458:LYS:N	2.34	0.41
1:A:182:ILE:HD12	1:A:258:VAL:HG22	2.02	0.41
1:A:740:ASP:N	1:A:740:ASP:OD1	2.51	0.41
1:A:328:ILE:HG22	1:A:335:VAL:CG1	2.51	0.41
1:A:741:MET:HE3	1:A:766:ILE:HD11	2.02	0.41
1:A:768:MET:HG2	1:A:792:LYS:HB3	2.02	0.41
1:A:479:THR:OG1	1:A:480:VAL:N	2.53	0.41
1:A:628:VAL:HG12	1:A:629:VAL:N	2.36	0.41
2:B:220:LEU:HD23	2:B:223:GLN:H	1.86	0.41
6:F:230:LEU:O	6:F:230:LEU:HD23	2.20	0.41
6:F:447:ILE:HG23	6:F:471:ILE:HD11	2.02	0.41
1:A:601:ILE:HD13	1:A:601:ILE:HA	1.98	0.41
3:C:218:ASN:OD1	3:C:222:ASN:N	2.54	0.41
1:A:727:ASP:OD1	1:A:728:LYS:N	2.54	0.41
1:A:606:ASN:O	1:A:607:GLU:C	2.64	0.40
1:A:197:PHE:CZ	1:A:226:LEU:HD12	2.55	0.40
2:B:227:ILE:HG22	2:B:228:TRP:HD1	1.86	0.40
6:F:475:ARG:O	6:F:475:ARG:HD3	2.21	0.40
1:A:494:PHE:CZ	1:A:496:ALA:HB2	2.57	0.40
2:B:253:VAL:HG22	2:B:258:VAL:HG22	2.04	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	A	768/790 (97%)	714 (93%)	50 (6%)	4 (0%)	24	57
2	B	361/373 (97%)	352 (98%)	9 (2%)	0	100	100
3	C	317/320 (99%)	297 (94%)	17 (5%)	3 (1%)	14	46
4	D	215/226 (95%)	211 (98%)	2 (1%)	2 (1%)	14	46
5	E	90/104 (86%)	88 (98%)	2 (2%)	0	100	100
6	F	457/492 (93%)	451 (99%)	5 (1%)	1 (0%)	43	72
All	All	2208/2305 (96%)	2113 (96%)	85 (4%)	10 (0%)	26	57

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	685	ASP
1	A	686	TYR
3	C	227	THR
1	A	680	SER
4	D	207	ASP
6	F	30	ALA
1	A	447	ASP
3	C	219	ALA
3	C	311	ASP
4	D	206	PRO

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	640/672 (95%)	602 (94%)	38 (6%)	18	45
2	B	295/304 (97%)	285 (97%)	10 (3%)	32	57
3	C	216/258 (84%)	213 (99%)	3 (1%)	59	71
4	D	182/190 (96%)	178 (98%)	4 (2%)	45	65
5	E	81/90 (90%)	81 (100%)	0	100	100
6	F	371/395 (94%)	368 (99%)	3 (1%)	73	77
All	All	1785/1909 (94%)	1727 (97%)	58 (3%)	35	58

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	211	ASP
1	A	230	TYR
1	A	252	LYS
1	A	258	VAL
1	A	261	THR
1	A	322	VAL
1	A	323	GLN
1	A	324	SER
1	A	328	ILE
1	A	332	ASP
1	A	345	ASN
1	A	349	VAL
1	A	351	LYS
1	A	352	ILE
1	A	353	ARG
1	A	359	THR
1	A	399	ASP
1	A	422	ASN
1	A	537	SER
1	A	573	ASN
1	A	577	THR
1	A	581	LEU
1	A	605	ASP
1	A	614	ASP
1	A	615	THR
1	A	629	VAL
1	A	630	LEU
1	A	632	ARG
1	A	641	LEU

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Mol	Chain	Res	Type
1	A	650	GLU
1	A	690	CYS
1	A	728	LYS
1	A	766	ILE
1	A	769	SER
1	A	774	LEU
1	A	778	SER
1	A	801	GLN
2	B	23	PHE
2	B	27	GLU
2	B	29	VAL
2	B	31	LYS
2	B	32	MET
2	B	54	SER
2	B	88	ASP
2	B	111	LEU
2	B	220	LEU
2	B	323	LEU
3	C	100	THR
3	C	221	GLN
3	C	331	ASN
4	D	181	VAL
4	D	199	GLU
4	D	202	LEU
4	D	237	ILE
6	F	29	SER
6	F	140	HIS
6	F	198	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	179	GLN
1	A	198	GLN
1	A	286	GLN
1	A	323	GLN
1	A	441	GLN
1	A	594	ASN
2	B	175	ASN
2	B	301	GLN
3	C	114	ASN

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Mol	Chain	Res	Type
3	C	149	ASN
4	D	116	ASN
4	D	228	ASN
4	D	230	GLN
6	F	199	GLN
6	F	232	GLN
6	F	418	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

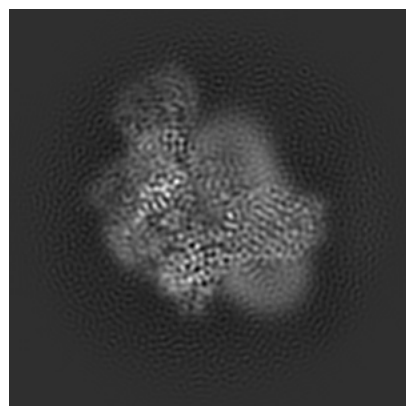
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-56543. These allow visual inspection of the internal detail of the map and identification of artifacts.

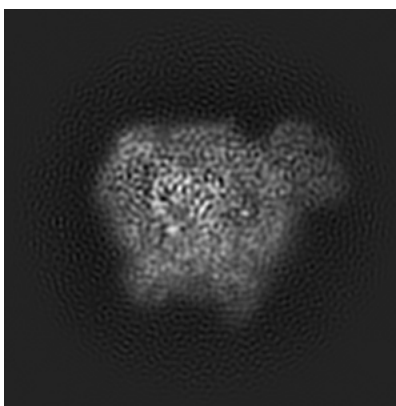
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

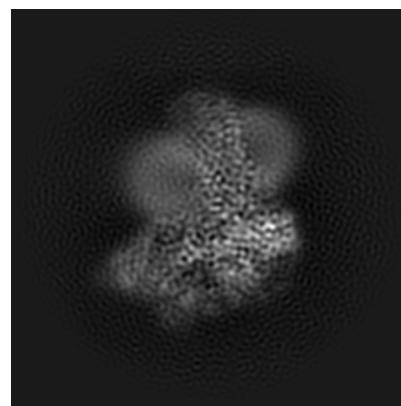
6.1.1 Primary map



X

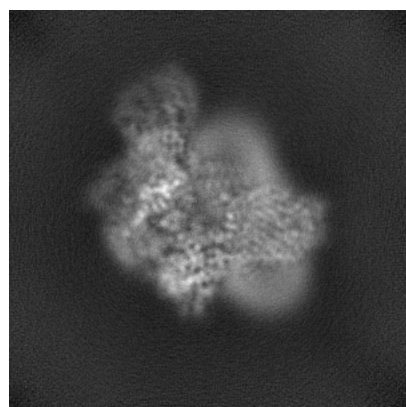


Y

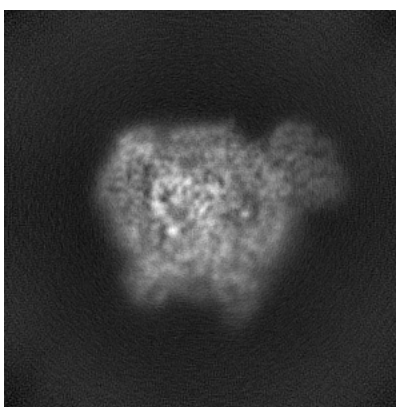


Z

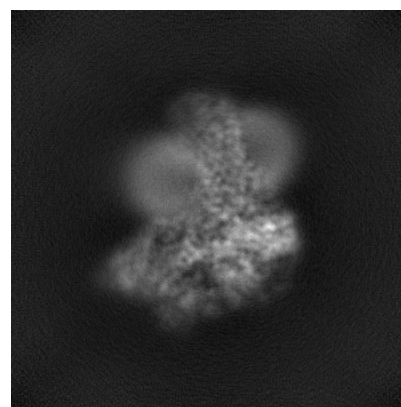
6.1.2 Raw map



X



Y

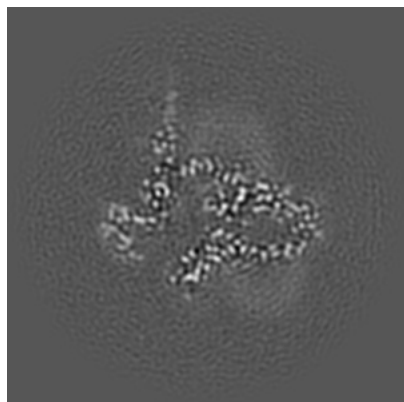


Z

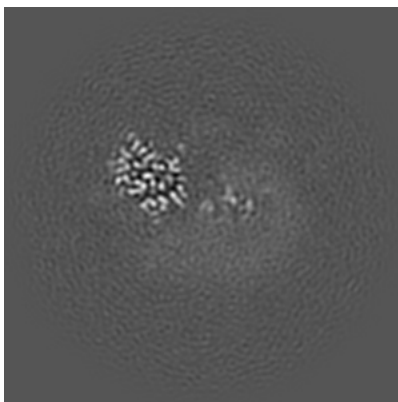
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

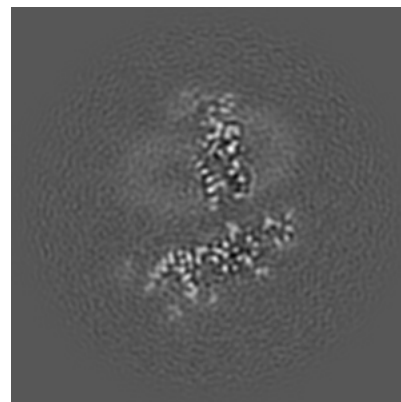
6.2.1 Primary map



X Index: 150

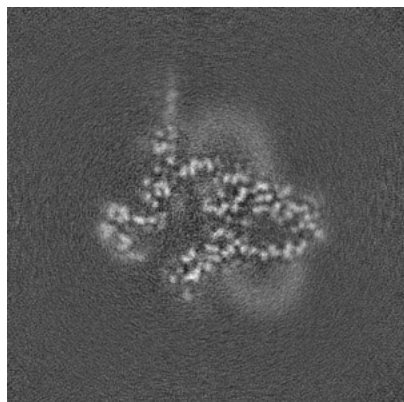


Y Index: 150

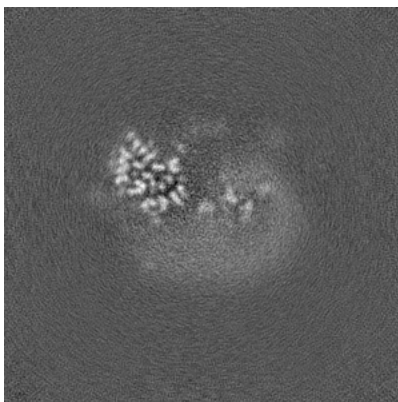


Z Index: 150

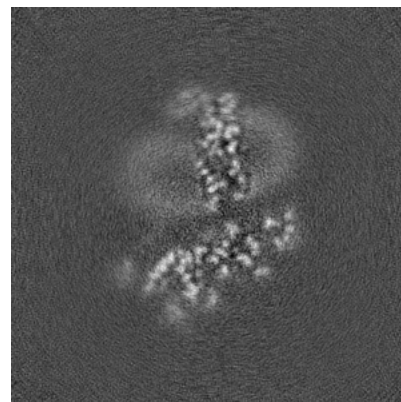
6.2.2 Raw map



X Index: 150



Y Index: 150

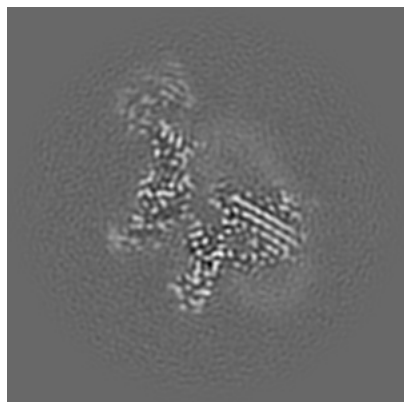


Z Index: 150

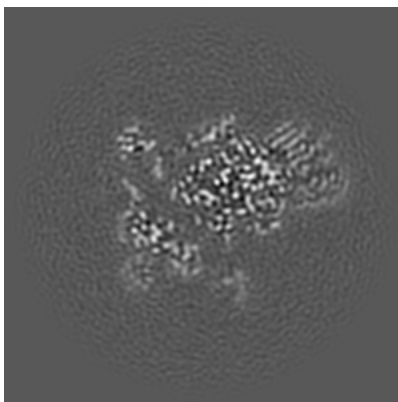
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

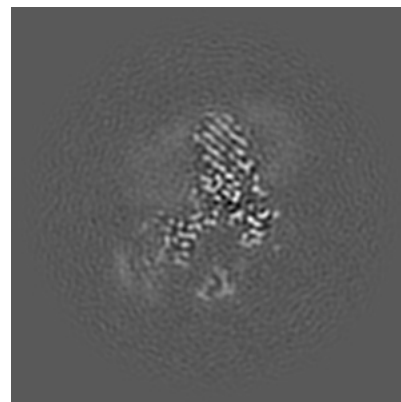
6.3.1 Primary map



X Index: 169

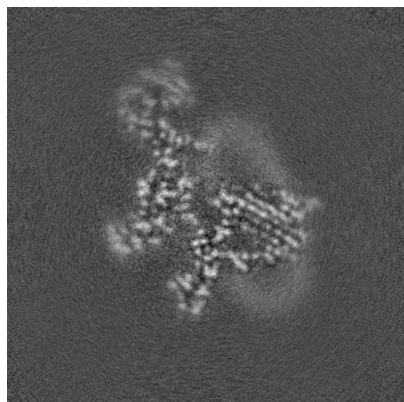


Y Index: 122

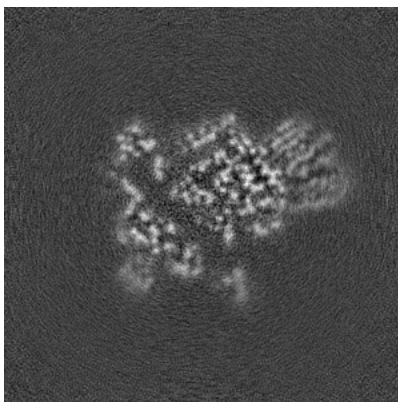


Z Index: 115

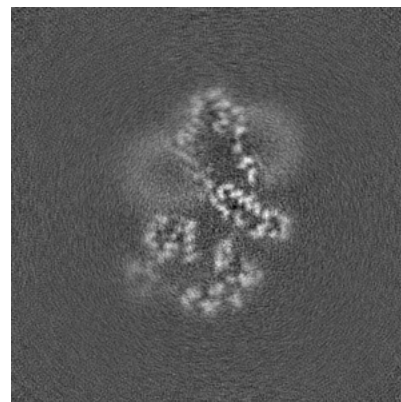
6.3.2 Raw map



X Index: 167



Y Index: 123

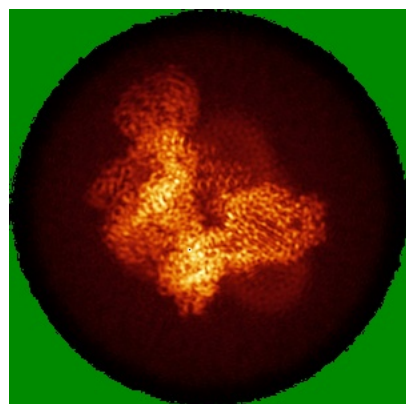


Z Index: 128

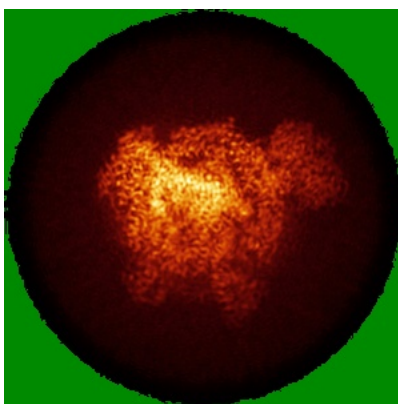
The images above show the largest variance slices of the map in three orthogonal directions.

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

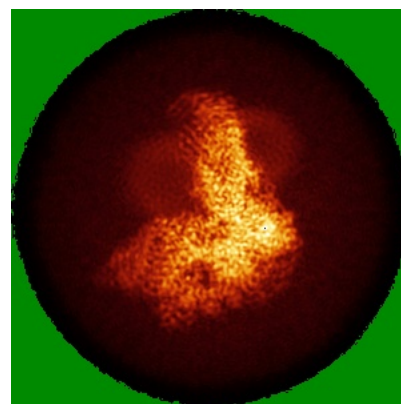
6.4.1 Primary map



X

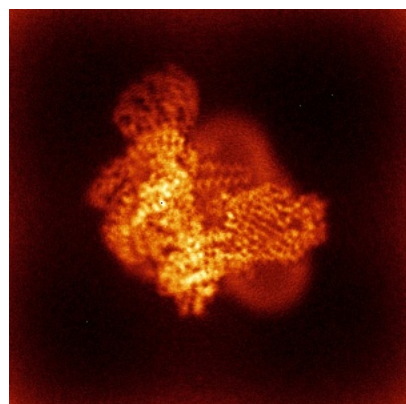


Y

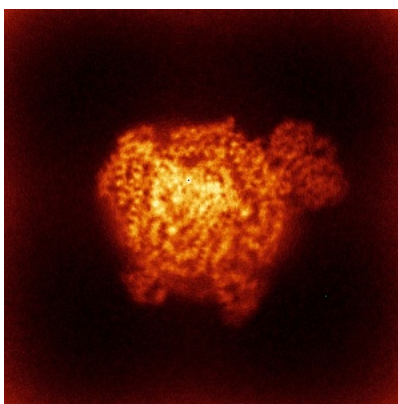


Z

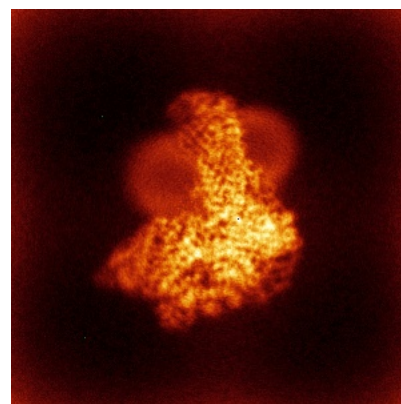
6.4.2 Raw map



X



Y

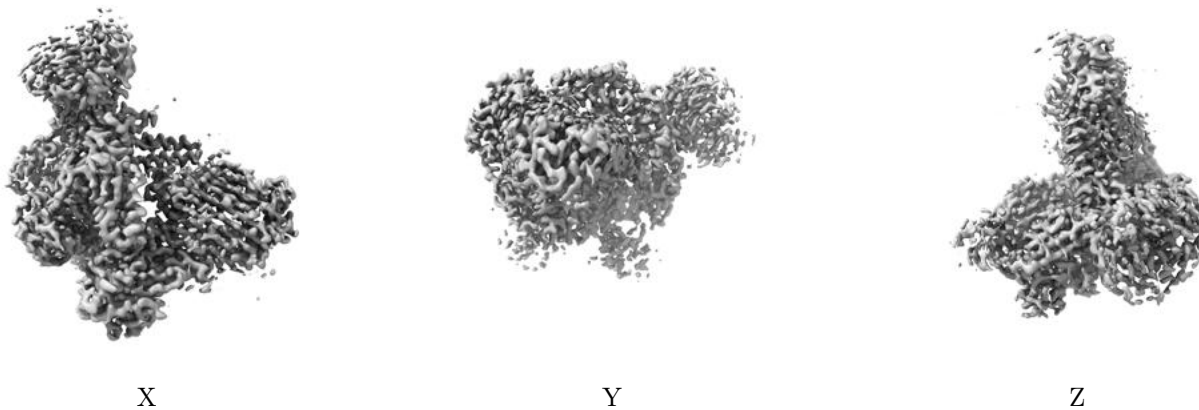


Z

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.

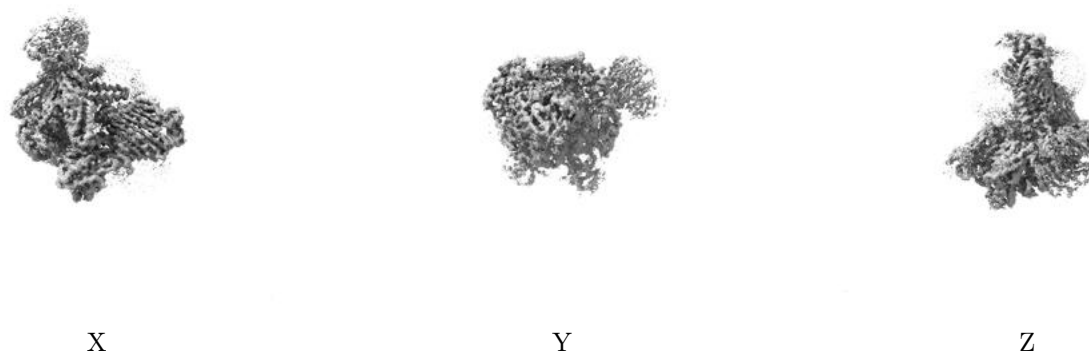
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.145. 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.

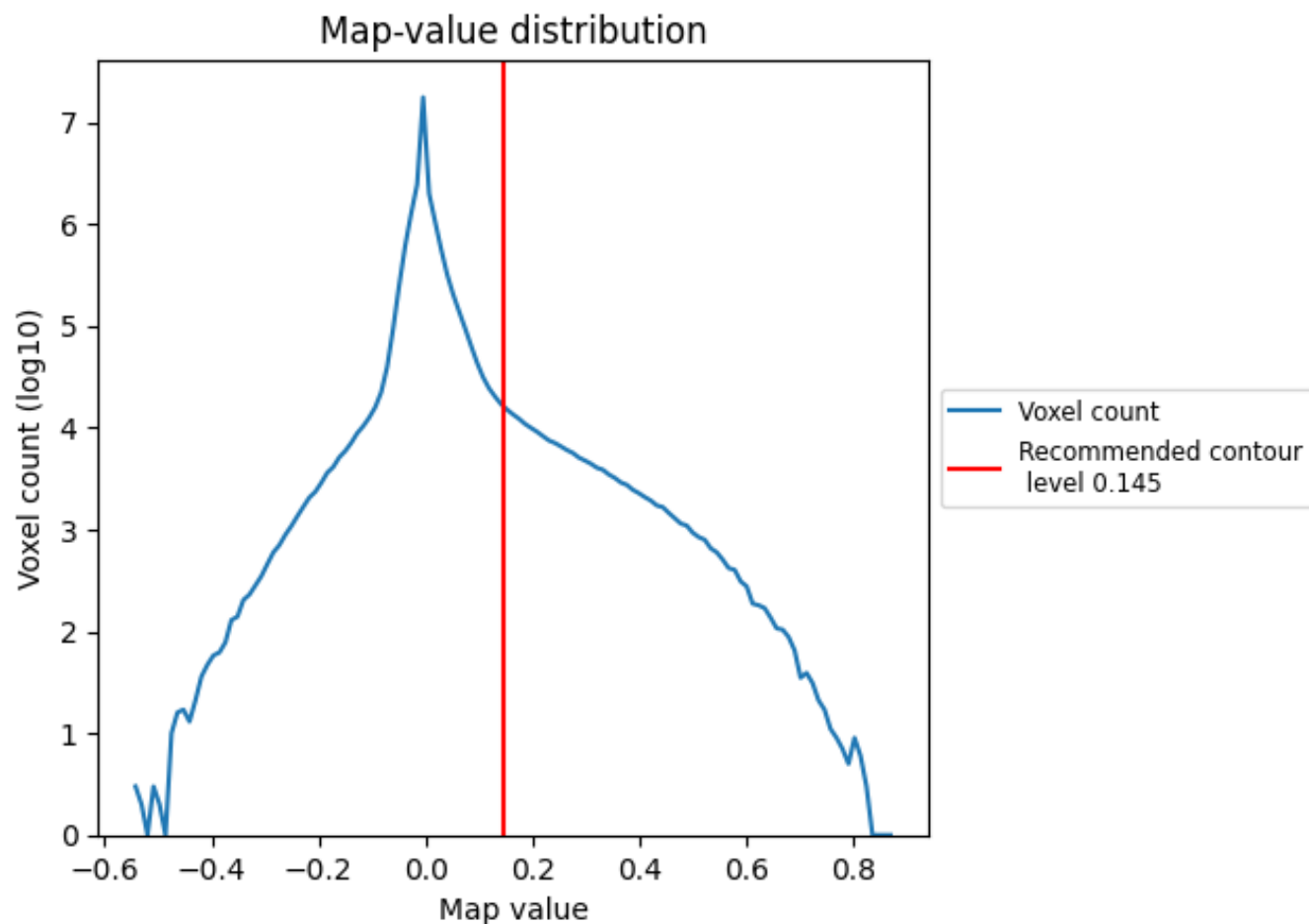
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

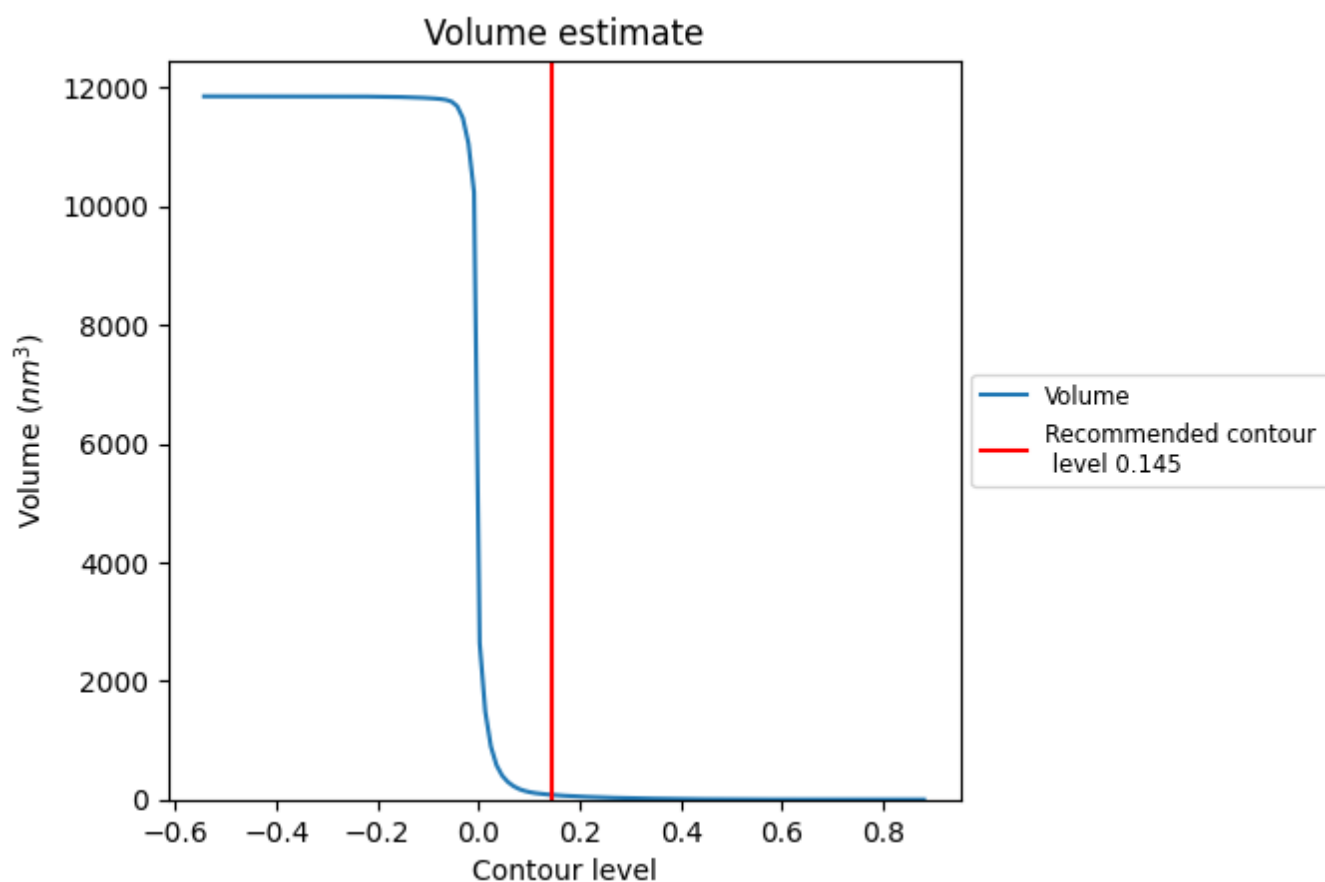
This section contains the results of statistical analysis of the map.

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.

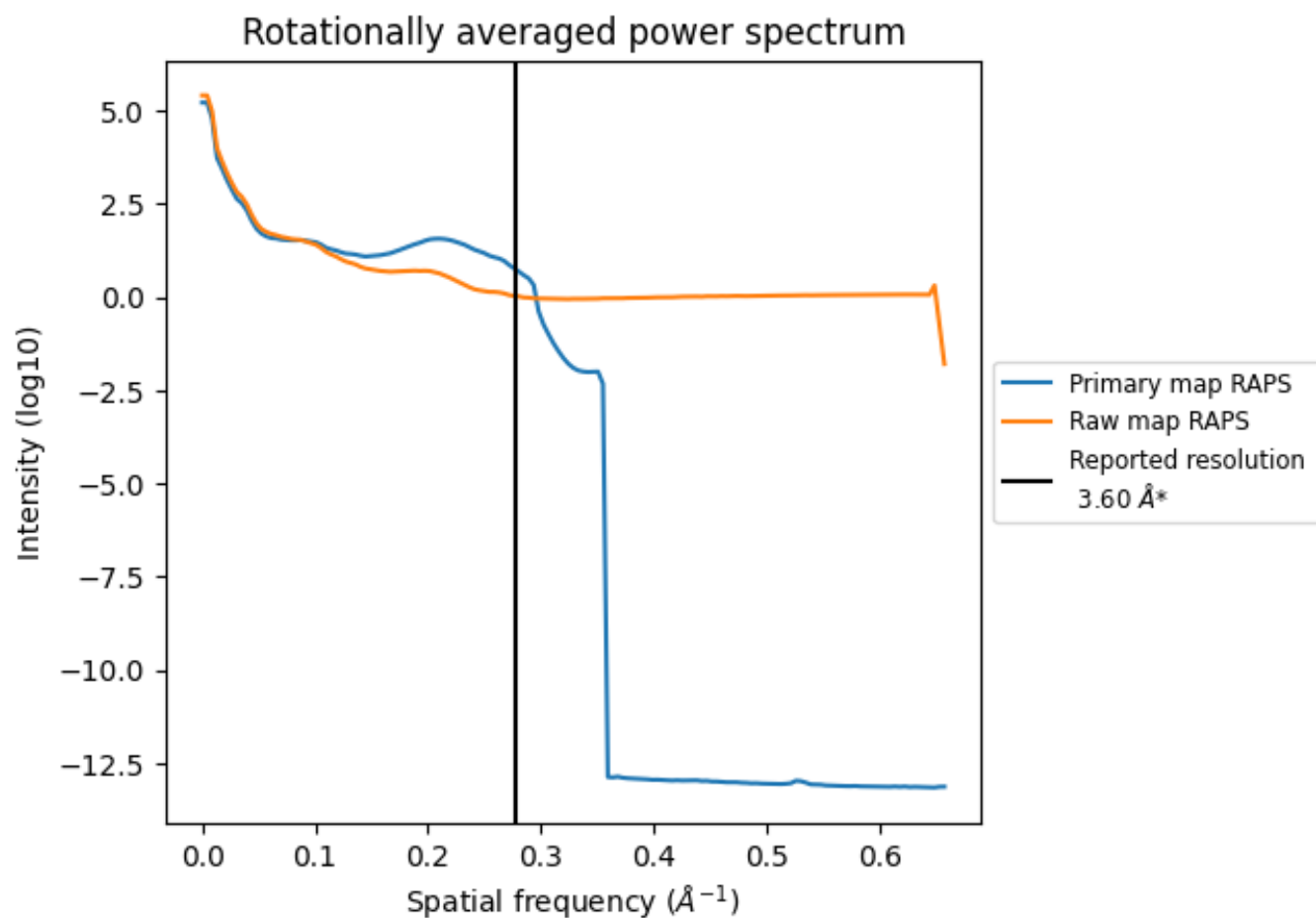
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 80 nm³; this corresponds to an approximate mass of 72 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 ⓘ

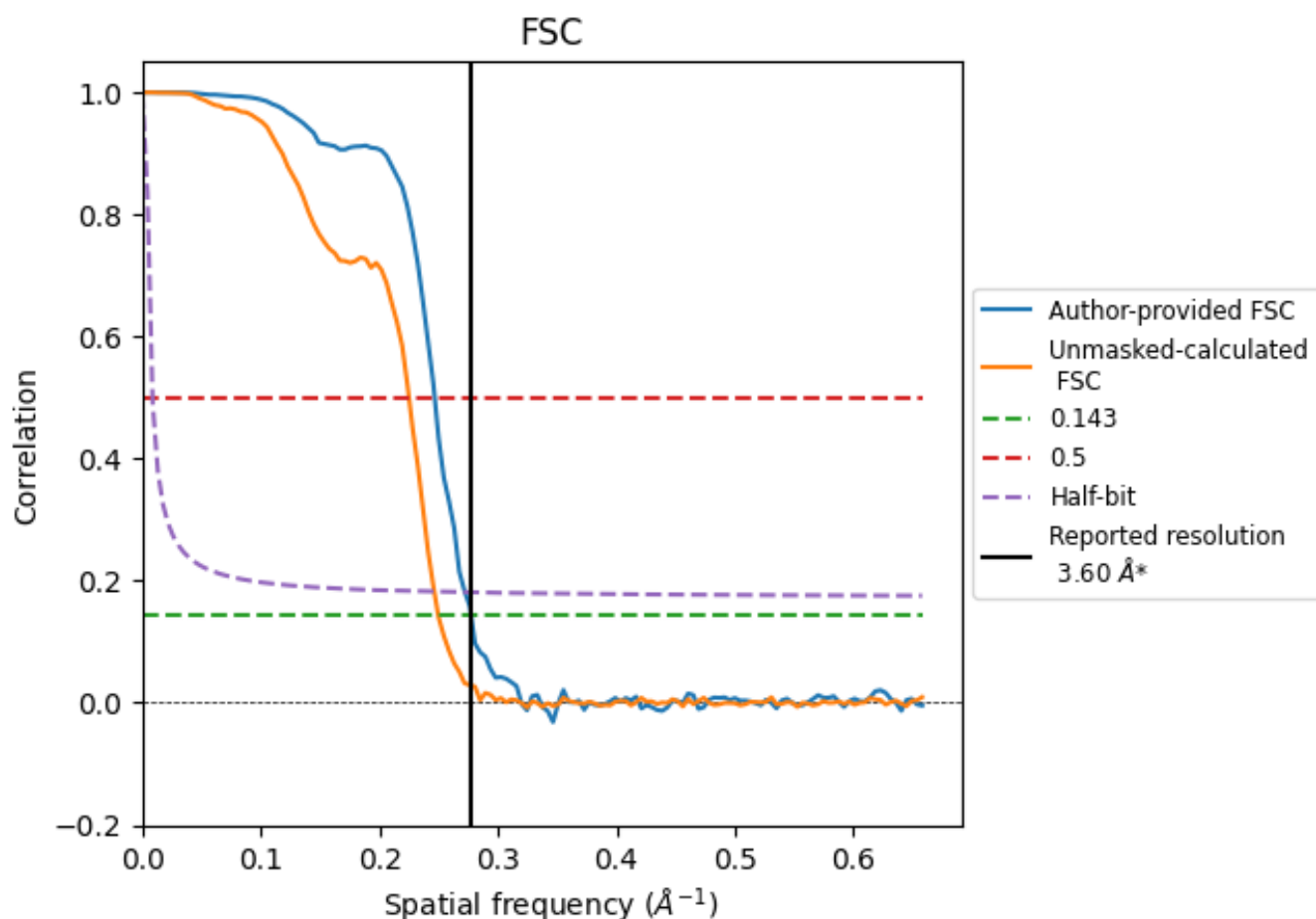


*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

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 0.278 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

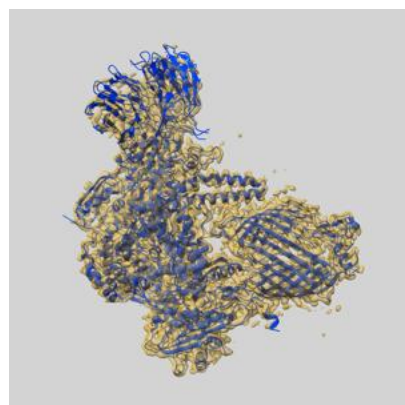
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.60	-	-
Author-provided FSC curve	3.60	4.05	3.67
Unmasked-calculated*	4.00	4.44	4.06

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.00 differs from the reported value 3.6 by more than 10 %

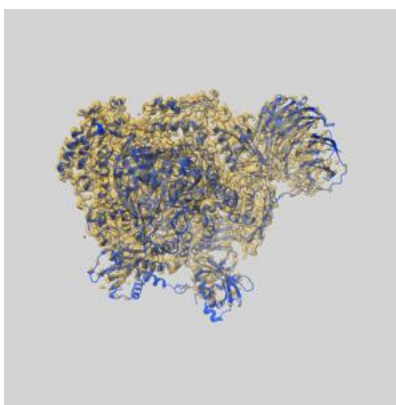
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-56543 and PDB model 28JL. Per-residue inclusion information can be found in section 3 on page 7.

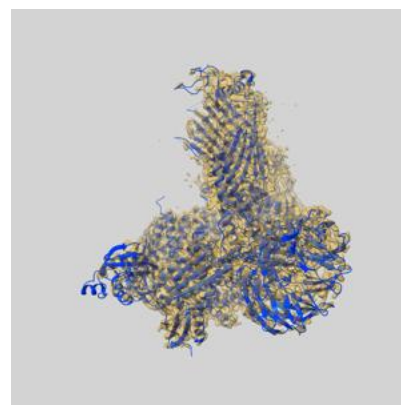
9.1 Map-model overlay [i](#)



X



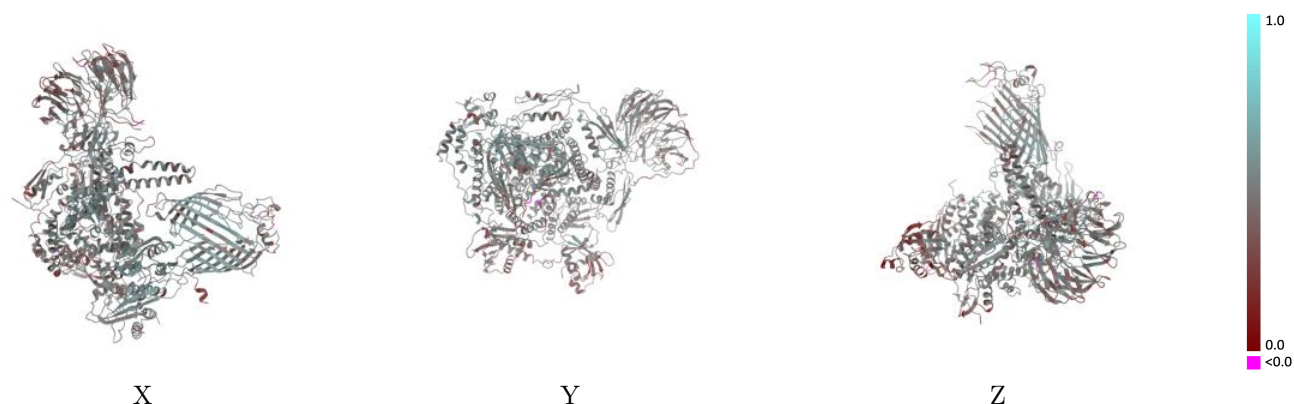
Y



Z

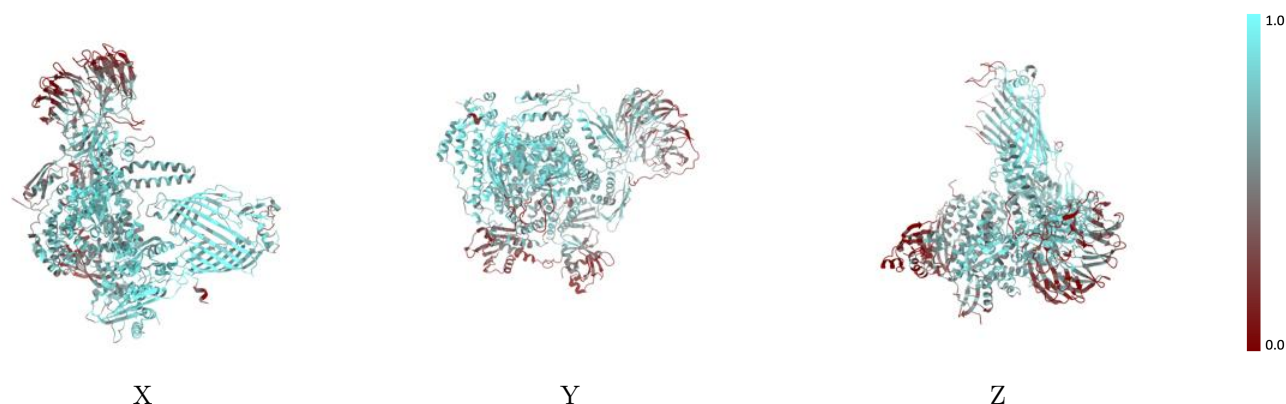
The images above show the 3D surface view of the map at the recommended contour level 0.145 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



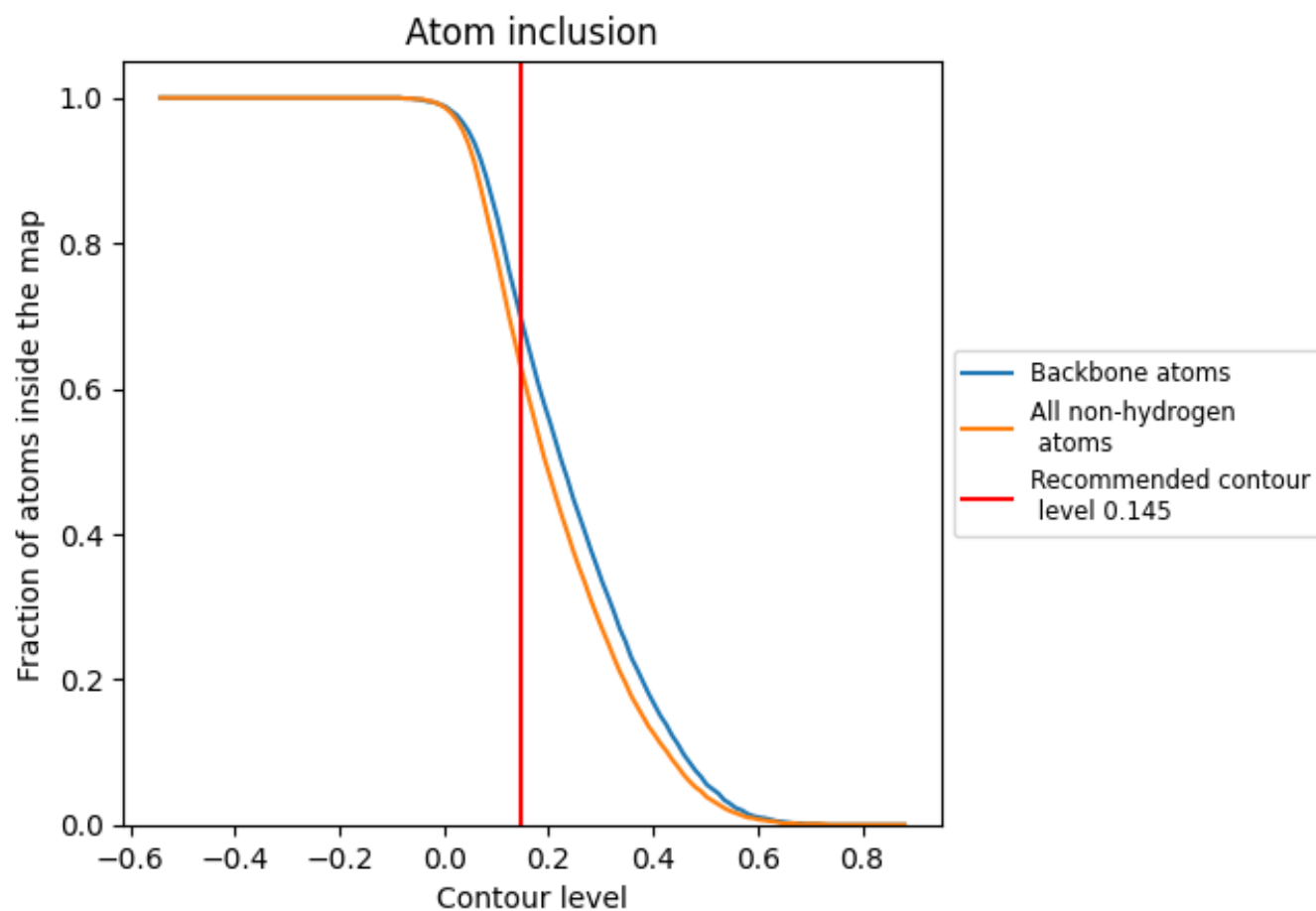
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.145).

9.4 Atom inclusion [i](#)



At the recommended contour level, 70% of all backbone atoms, 64% 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 (0.145) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6350	0.4580
A	0.7260	0.4760
B	0.4000	0.4200
C	0.3230	0.3750
D	0.7820	0.4900
E	0.7520	0.4820
F	0.7660	0.4900

1.0

0.0

<0.0