



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 11:43 AM UTC

PDB ID : 8QN4 / pdb_00008qn4
EMDB ID : EMD-18507
Title : Structure of the BAM-EspP complex in an open EspP intermediate state
Authors : Xie, T.; Pang, J.; Shen, C.; Chang, S.; Tang, X.; Zhang, X.; Dong, H.; Zhou, R.
Deposited on : 2023-09-25
Resolution : 3.36 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:

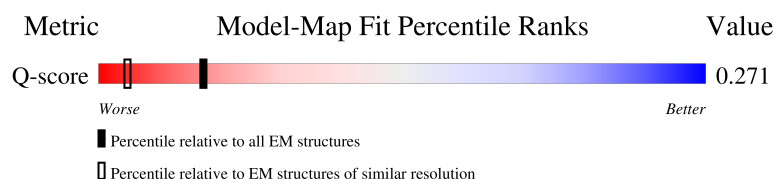
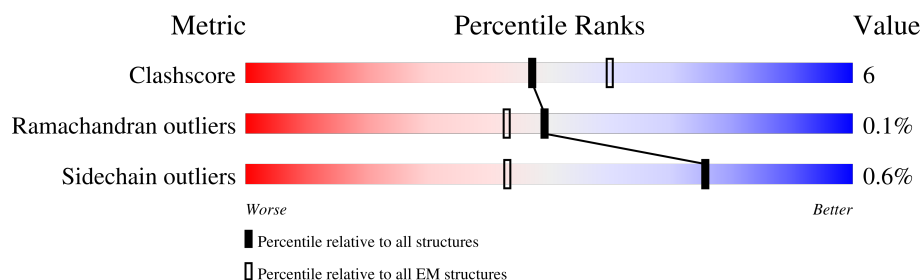
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.36 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	14332 (2.86 - 3.86)

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	810	
2	B	392	
3	C	344	
4	D	245	

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Mol	Chain	Length	Quality of chain
5	E	123	14% 59% 8% 33%
6	F	353	14% 58% 8% 33%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12443 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	636	Total	C	N	O	S	0	0
			5066	3199	847	1005	15		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	431	CYS	GLY	engineered mutation	UNP P0A940

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	367	Total	C	N	O	S	0	0
			2750	1725	470	549	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	56	Total	C	N	O	S	0	0
			407	256	71	79	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	213	Total	C	N	O	S	0	0
			1722	1086	303	326	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	83	Total	C	N	O	S	0	0
			645	405	112	126	2		

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

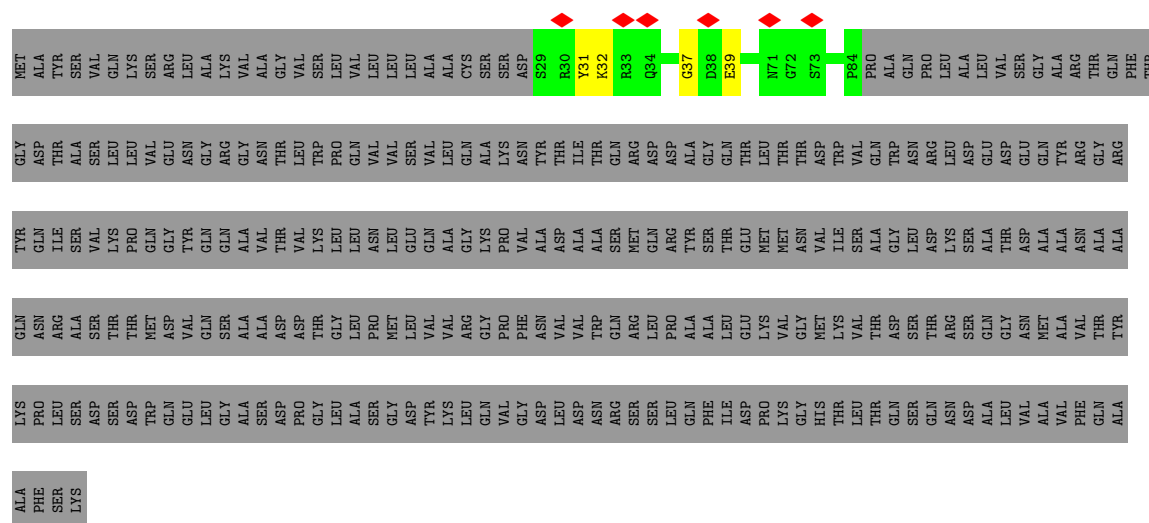
- Molecule 6 is a protein called EspP epsilon.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	236	Total	C	N	O	S	0	0
			1853	1168	319	360	6		

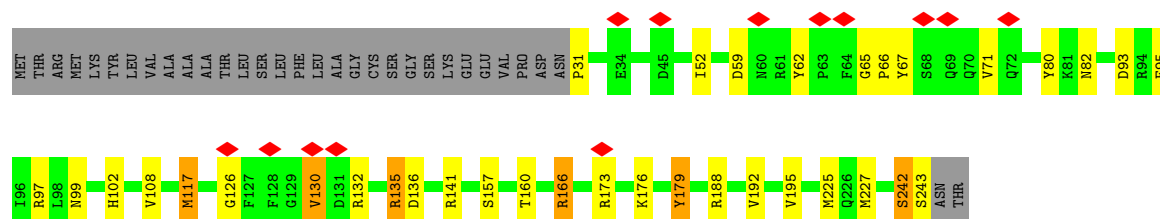
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	1293	CYS	ASN	engineered mutation	UNP C1J8F9

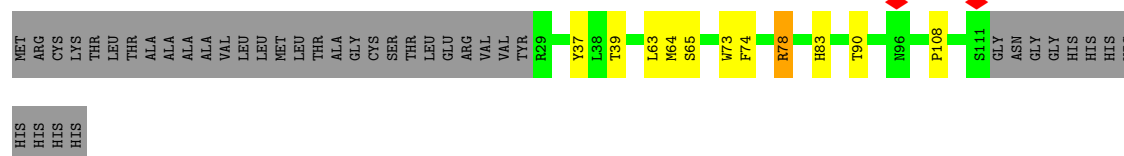
Chain C: 15% 84%



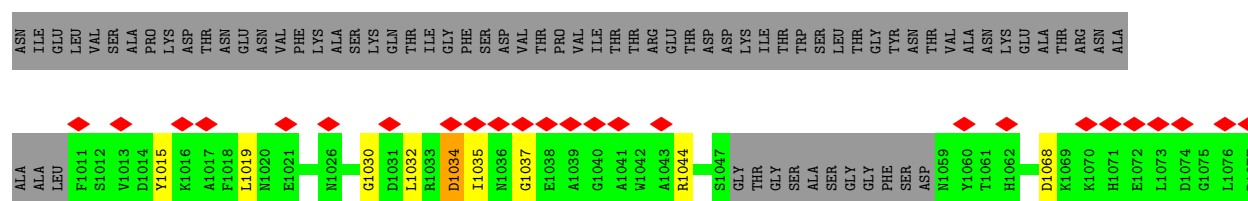
Chain D: 5% 72% 12% 13%

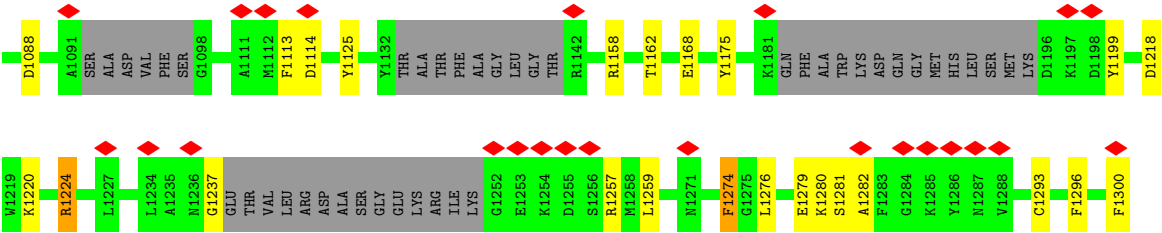


Chain E: 59% 8% 33%



Chain F: 14% 58% 8% 33%





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	162981	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.402	Depositor
Minimum map value	-0.782	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.028	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	238.08, 238.08, 238.08	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.93, 0.93, 0.93	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1.02	0/5196	1.46	33/7053 (0.5%)
2	B	1.03	2/2800 (0.1%)	1.43	11/3819 (0.3%)
3	C	0.99	0/415	1.34	0/565
4	D	1.01	0/1761	1.55	10/2390 (0.4%)
5	E	1.02	0/659	1.37	4/899 (0.4%)
6	F	1.09	0/1891	1.41	13/2542 (0.5%)
All	All	1.03	2/12722 (0.0%)	1.45	71/17268 (0.4%)

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	9
2	B	0	3
4	D	0	4
5	E	0	1
6	F	0	5
All	All	0	22

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	286	VAL	CA-C	-5.63	1.47	1.52
2	B	210	GLY	CA-C	-5.21	1.47	1.52

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	196	GLY	N-CA-C	-17.68	94.01	112.04
2	B	286	VAL	CB-CA-C	-9.28	102.79	111.05
1	A	493	ASP	CA-CB-CG	8.62	121.22	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	287	ASN	CA-CB-CG	-8.45	104.15	112.60
4	D	242	SER	N-CA-C	8.16	128.19	110.80
1	A	432	TYR	N-CA-C	-8.14	96.63	109.50
1	A	385	GLY	CA-C-N	7.89	130.85	120.28
1	A	385	GLY	C-N-CA	7.89	130.85	120.28
6	F	1034	ASP	N-CA-CB	-7.87	98.47	110.91
4	D	243	SER	N-CA-CB	6.94	122.30	110.50
1	A	277	LEU	N-CA-C	-6.80	105.27	113.97
1	A	498	ASP	CA-CB-CG	6.71	119.31	112.60
1	A	660	VAL	N-CA-C	-6.64	95.69	107.18
1	A	657	SER	N-CA-C	6.58	119.04	111.02
1	A	401	ASP	CA-CB-CG	6.57	119.17	112.60
2	B	310	ILE	N-CA-C	6.56	117.31	110.62
6	F	1113	PHE	CA-C-N	6.50	128.99	120.28
6	F	1113	PHE	C-N-CA	6.50	128.99	120.28
2	B	50	THR	N-CA-CB	6.38	119.54	109.71
1	A	310	LYS	N-CA-C	6.22	118.14	111.36
4	D	141	ARG	CA-C-N	6.09	128.45	120.28
4	D	141	ARG	C-N-CA	6.09	128.45	120.28
6	F	1114	ASP	CA-C-N	6.00	128.24	120.44
6	F	1114	ASP	C-N-CA	6.00	128.24	120.44
1	A	708	GLY	N-CA-C	-5.91	104.66	112.81
6	F	1162	THR	N-CA-C	-5.88	100.37	109.14
5	E	39	THR	CA-C-N	5.86	128.14	120.28
5	E	39	THR	C-N-CA	5.86	128.14	120.28
4	D	135	ARG	NE-CZ-NH2	5.81	124.43	119.20
1	A	239	ASN	CA-CB-CG	5.78	118.38	112.60
1	A	407	GLY	N-CA-C	-5.78	102.62	110.43
1	A	580	LYS	N-CA-C	-5.78	104.46	112.03
1	A	709	ASN	CA-CB-CG	5.72	118.32	112.60
6	F	1030	GLY	CA-C-N	5.72	127.94	120.28
6	F	1030	GLY	C-N-CA	5.72	127.94	120.28
1	A	662	GLY	N-CA-C	-5.62	107.62	114.92
4	D	65	GLY	CA-C-N	5.58	126.82	119.84
4	D	65	GLY	C-N-CA	5.58	126.82	119.84
1	A	525	LEU	N-CA-C	-5.55	101.13	109.95
5	E	83	HIS	N-CA-C	-5.46	106.54	113.43
1	A	212	ARG	NE-CZ-NH2	5.45	124.11	119.20
6	F	1218	ASP	N-CA-CB	-5.45	103.35	110.88
1	A	307	ASP	CA-C-N	5.45	127.92	120.46
1	A	307	ASP	C-N-CA	5.45	127.92	120.46
2	B	240	GLU	N-CA-C	-5.42	107.68	114.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	588	THR	N-CA-C	-5.41	107.33	114.31
6	F	1015	TYR	CA-C-N	5.36	127.47	120.28
6	F	1015	TYR	C-N-CA	5.36	127.47	120.28
1	A	734	ARG	NE-CZ-NH2	5.34	124.00	119.20
2	B	246	ASP	CA-CB-CG	-5.33	107.28	112.60
1	A	336	LYS	N-CA-C	-5.32	101.09	109.50
6	F	1068	ASP	CA-CB-CG	5.31	117.91	112.60
1	A	464	ASP	CA-C-N	5.22	127.73	120.63
1	A	464	ASP	C-N-CA	5.22	127.73	120.63
1	A	669	GLY	CA-C-N	5.21	126.35	119.84
1	A	669	GLY	C-N-CA	5.21	126.35	119.84
1	A	725	ILE	N-CA-C	-5.20	106.82	113.22
5	E	78	ARG	NE-CZ-NH2	5.19	123.88	119.20
4	D	188	ARG	N-CA-C	-5.15	107.05	113.38
4	D	117	MET	CG-SD-CE	-5.15	89.58	100.90
2	B	196	GLY	CA-C-N	-5.14	112.90	121.94
2	B	196	GLY	C-N-CA	-5.14	112.90	121.94
1	A	639	ASP	CA-C-N	5.08	125.98	120.44
1	A	639	ASP	C-N-CA	5.08	125.98	120.44
1	A	400	THR	CA-C-N	5.07	129.92	122.77
1	A	400	THR	C-N-CA	5.07	129.92	122.77
6	F	1088	ASP	CA-CB-CG	5.05	117.65	112.60
4	D	108	VAL	N-CA-C	5.03	115.26	110.53
1	A	661	ARG	NE-CZ-NH2	5.02	123.72	119.20
2	B	363	ASP	CA-C-N	5.02	127.01	120.28
2	B	363	ASP	C-N-CA	5.02	127.01	120.28

There are no chirality outliers.

All (22) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	229	TYR	Sidechain
1	A	230	TYR	Sidechain
1	A	255	TYR	Sidechain
1	A	391	ARG	Sidechain
1	A	465	TYR	Sidechain
1	A	504	TYR	Sidechain
1	A	686	TYR	Sidechain
1	A	688	TYR	Sidechain
1	A	789	GLN	Peptide
2	B	280	LYS	Peptide
2	B	355	ARG	Sidechain

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Mol	Chain	Res	Type	Group
2	B	64	HIS	Sidechain
4	D	166	ARG	Sidechain
4	D	173	ARG	Sidechain
4	D	179	TYR	Sidechain
4	D	80	TYR	Sidechain
5	E	78	ARG	Sidechain
6	F	1044	ARG	Sidechain
6	F	1125	TYR	Sidechain
6	F	1175	TYR	Sidechain
6	F	1224	ARG	Sidechain
6	F	1274	PHE	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5066	0	4756	33	0
2	B	2750	0	2682	61	0
3	C	407	0	407	3	0
4	D	1722	0	1668	31	0
5	E	645	0	625	13	0
6	F	1853	0	1740	24	0
All	All	12443	0	11878	147	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 6.

All (147) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:258:VAL:HG13	2:B:272:LEU:HD11	1.42	0.99
2:B:322:LEU:HD21	2:B:327:LEU:HD11	1.42	0.99
2:B:283:LEU:CD1	2:B:308:LEU:HD11	2.03	0.89
2:B:202:THR:HG22	2:B:221:MET:HE1	1.62	0.82
1:A:348:TYR:HB3	5:E:63:LEU:HD22	1.63	0.80
1:A:287:LEU:HD22	1:A:311:LEU:HD11	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:192:VAL:HG13	5:E:64:MET:HE3	1.63	0.79
1:A:657:SER:HA	1:A:661:ARG:CG	2.16	0.76
2:B:64:HIS:CE1	2:B:154:ARG:HD3	2.21	0.75
2:B:35:LEU:HD21	2:B:325:ARG:HH11	1.53	0.72
4:D:67:TYR:CE1	4:D:71:VAL:HG22	2.25	0.72
6:F:1259:LEU:HD11	6:F:1279:GLU:CD	2.15	0.72
1:A:283:GLU:O	1:A:287:LEU:HD13	1.92	0.70
1:A:266:TYR:HB3	1:A:335:VAL:HG23	1.74	0.69
1:A:657:SER:HA	1:A:661:ARG:CD	2.24	0.68
4:D:117:MET:HE1	4:D:166:ARG:NH1	2.09	0.67
2:B:152:LEU:HD11	2:B:195:ARG:NH1	2.10	0.67
4:D:227:MET:HE2	5:E:90:THR:HG21	1.76	0.66
6:F:1257:ARG:HE	6:F:1281:SER:HB3	1.61	0.65
1:A:411:GLN:HE22	5:E:63:LEU:HD13	1.62	0.65
2:B:327:LEU:HD12	2:B:339:VAL:HB	1.79	0.64
4:D:117:MET:HE1	4:D:166:ARG:HH11	1.63	0.64
4:D:62:TYR:CD1	4:D:66:PRO:HD2	2.33	0.64
2:B:283:LEU:HD12	2:B:308:LEU:HD11	1.79	0.64
2:B:325:ARG:O	2:B:327:LEU:HD22	1.98	0.63
4:D:192:VAL:HG13	5:E:64:MET:CE	2.28	0.63
2:B:194:LEU:HG	2:B:195:ARG:H	1.64	0.63
2:B:35:LEU:HD21	2:B:325:ARG:NH1	2.13	0.62
2:B:158:SER:CB	2:B:202:THR:HG21	2.30	0.62
1:A:686:TYR:CD2	1:A:688:TYR:CE2	2.88	0.61
6:F:1259:LEU:HD21	6:F:1279:GLU:OE2	1.99	0.61
6:F:1274:PHE:CE1	6:F:1296:PHE:CE1	2.89	0.61
6:F:1035:ILE:HG21	6:F:1220:LYS:HE2	1.83	0.61
1:A:686:TYR:CE2	1:A:688:TYR:CZ	2.89	0.60
2:B:90:LYS:HG2	2:B:91:GLU:H	1.67	0.59
2:B:48:TRP:CZ3	2:B:84:LEU:HD13	2.38	0.58
2:B:323:LEU:HG	2:B:324:HIS:CD2	2.38	0.58
4:D:192:VAL:HG22	4:D:225:MET:HE1	1.85	0.58
1:A:657:SER:HA	1:A:661:ARG:HD3	1.85	0.57
2:B:270:LEU:HD22	2:B:275:GLY:O	2.05	0.56
4:D:227:MET:HE2	5:E:90:THR:CG2	2.35	0.56
1:A:657:SER:HA	1:A:661:ARG:HG3	1.85	0.56
2:B:347:HIS:CE1	2:B:359:GLN:CD	2.83	0.56
2:B:189:MET:SD	2:B:212:ASP:HB3	2.46	0.55
2:B:270:LEU:CD2	2:B:278:MET:HB3	2.38	0.54
6:F:1019:LEU:CD2	6:F:1257:ARG:HH12	2.21	0.54
2:B:158:SER:HB2	2:B:202:THR:HG21	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:290:ILE:HG12	2:B:330:PRO:O	2.08	0.53
5:E:65:SER:HG	5:E:73:TRP:CD1	2.26	0.53
2:B:263:TYR:CD2	2:B:286:VAL:CG2	2.92	0.52
2:B:303:ASP:HB3	2:B:327:LEU:HD23	1.90	0.52
4:D:157:SER:O	4:D:160:THR:HG22	2.09	0.52
6:F:1035:ILE:HG22	6:F:1035:ILE:O	2.09	0.52
4:D:67:TYR:CE1	4:D:71:VAL:CG2	2.91	0.52
2:B:131:VAL:HG11	2:B:164:ILE:CD1	2.40	0.52
2:B:195:ARG:HG2	2:B:196:GLY:H	1.75	0.52
6:F:1019:LEU:HD21	6:F:1257:ARG:HH12	1.74	0.52
6:F:1034:ASP:OD1	6:F:1224:ARG:HD3	2.10	0.51
4:D:99:ASN:HD22	4:D:102:HIS:CD2	2.28	0.51
6:F:1199:TYR:HB3	6:F:1237:GLY:H	1.75	0.51
2:B:195:ARG:CG	2:B:196:GLY:H	2.23	0.50
2:B:281:ARG:HH12	2:B:283:LEU:HD21	1.76	0.50
1:A:188:PHE:HE2	1:A:193:LEU:HG	1.76	0.50
1:A:767:ARG:HD3	1:A:794:TYR:HB2	1.93	0.50
4:D:52:ILE:HG21	4:D:82:ASN:ND2	2.27	0.50
2:B:48:TRP:CH2	2:B:84:LEU:HD13	2.47	0.50
6:F:1280:LYS:HE2	6:F:1282:ALA:HB2	1.94	0.50
2:B:345:TYR:CE2	2:B:361:LYS:HE3	2.47	0.49
6:F:1034:ASP:HB3	6:F:1224:ARG:NH1	2.27	0.49
2:B:194:LEU:HD21	2:B:263:TYR:CD2	2.48	0.49
2:B:231:ARG:HH22	2:B:241:ILE:HG23	1.77	0.49
4:D:126:GLY:HA2	6:F:1300:PHE:C	2.37	0.49
2:B:131:VAL:HG11	2:B:164:ILE:HD11	1.94	0.48
1:A:498:ASP:CG	1:A:499:ALA:H	2.21	0.48
2:B:48:TRP:CZ3	2:B:387:VAL:O	2.66	0.48
3:C:32:LYS:HE3	3:C:39:GLU:HG3	1.96	0.48
2:B:302:ASN:HB3	2:B:324:HIS:CE1	2.48	0.48
6:F:1019:LEU:HD21	6:F:1257:ARG:NH1	2.29	0.48
1:A:709:ASN:HD21	1:A:747:THR:HG22	1.78	0.48
2:B:194:LEU:CD2	2:B:263:TYR:CD2	2.97	0.48
2:B:263:TYR:CD2	2:B:286:VAL:HG22	2.49	0.48
4:D:130:VAL:HG22	4:D:132:ARG:H	1.79	0.47
2:B:345:TYR:CD2	2:B:361:LYS:HE3	2.50	0.47
4:D:227:MET:HE1	5:E:74:PHE:HB3	1.97	0.47
1:A:737:PHE:CE1	1:A:772:ILE:HG12	2.49	0.47
2:B:35:LEU:CD2	2:B:325:ARG:HH11	2.24	0.47
4:D:67:TYR:CD1	4:D:71:VAL:HG22	2.49	0.47
4:D:95:PHE:CZ	4:D:102:HIS:CD2	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:263:TYR:HD2	2:B:286:VAL:HG22	1.80	0.47
2:B:287:ASN:HD21	2:B:329:SER:N	2.12	0.46
4:D:52:ILE:HG21	4:D:82:ASN:HD22	1.80	0.46
2:B:270:LEU:HD21	2:B:278:MET:HB3	1.97	0.46
1:A:361:LYS:HB3	4:D:135:ARG:NH1	2.31	0.46
2:B:345:TYR:CD1	2:B:361:LYS:HG3	2.50	0.46
1:A:348:TYR:CB	5:E:63:LEU:HD22	2.39	0.46
2:B:194:LEU:HG	2:B:195:ARG:N	2.30	0.46
6:F:1034:ASP:CG	6:F:1224:ARG:HD3	2.41	0.46
1:A:424:GLY:HA2	1:A:446:GLN:HA	1.99	0.45
6:F:1274:PHE:CD1	6:F:1296:PHE:CD1	3.04	0.45
2:B:347:HIS:CE1	2:B:359:GLN:NE2	2.85	0.45
6:F:1032:LEU:HA	6:F:1037:GLY:HA3	1.98	0.45
1:A:374:GLY:HA2	5:E:63:LEU:HD21	1.98	0.45
4:D:225:MET:HE1	5:E:64:MET:HE1	1.99	0.45
2:B:63:LEU:O	2:B:369:THR:HG23	2.16	0.45
4:D:59:ASP:HA	4:D:67:TYR:OH	2.16	0.45
2:B:203:ALA:HB1	2:B:272:LEU:HD21	1.98	0.45
4:D:67:TYR:CD1	4:D:67:TYR:C	2.95	0.45
2:B:131:VAL:HG21	2:B:164:ILE:CD1	2.47	0.45
4:D:99:ASN:ND2	4:D:102:HIS:CD2	2.85	0.44
6:F:1257:ARG:HE	6:F:1281:SER:CB	2.30	0.44
2:B:290:ILE:O	2:B:290:ILE:HG13	2.17	0.44
4:D:195:VAL:CG2	5:E:64:MET:HE2	2.47	0.44
2:B:290:ILE:HD13	2:B:331:VAL:HA	2.00	0.43
4:D:176:LYS:HA	4:D:179:TYR:HB3	2.00	0.43
6:F:1257:ARG:HD2	6:F:1281:SER:O	2.18	0.43
1:A:473:VAL:HG13	1:A:473:VAL:O	2.18	0.43
1:A:266:TYR:CB	1:A:335:VAL:HG23	2.46	0.43
1:A:686:TYR:CD2	1:A:686:TYR:C	2.97	0.43
2:B:48:TRP:HE3	2:B:50:THR:HG23	1.83	0.43
3:C:32:LYS:HE3	3:C:39:GLU:CG	2.48	0.43
1:A:247:LEU:O	2:B:194:LEU:HD12	2.19	0.42
1:A:361:LYS:HE3	1:A:452:THR:HG22	2.00	0.42
6:F:1276:LEU:HD12	6:F:1293:CYS:O	2.20	0.42
1:A:699:LEU:HD13	1:A:753:GLN:OE1	2.20	0.42
2:B:290:ILE:CD1	2:B:331:VAL:HA	2.50	0.42
1:A:255:TYR:CD2	2:B:195:ARG:NH2	2.88	0.42
2:B:90:LYS:HG2	2:B:91:GLU:N	2.34	0.42
2:B:239:THR:HG22	2:B:239:THR:O	2.20	0.41
6:F:1158:ARG:HH11	6:F:1168:GLU:CD	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:131:VAL:HG21	2:B:164:ILE:HD13	2.02	0.41
6:F:1279:GLU:OE1	6:F:1279:GLU:HA	2.20	0.41
1:A:360:SER:O	4:D:135:ARG:HD3	2.20	0.41
2:B:195:ARG:HG2	2:B:196:GLY:N	2.35	0.41
2:B:308:LEU:HD23	2:B:315:THR:HA	2.03	0.41
3:C:31:TYR:HB2	3:C:37:GLY:H	1.85	0.41
4:D:95:PHE:CE2	4:D:102:HIS:CD2	3.09	0.41
1:A:246:SER:HB3	2:B:195:ARG:HB2	2.01	0.41
1:A:767:ARG:CD	1:A:794:TYR:HB2	2.50	0.41
4:D:62:TYR:O	4:D:67:TYR:CZ	2.74	0.41
6:F:1259:LEU:HD11	6:F:1279:GLU:OE2	2.20	0.41
2:B:36:PRO:HG2	2:B:347:HIS:HE1	1.85	0.41
1:A:686:TYR:CD2	1:A:688:TYR:CD2	3.09	0.40
6:F:1274:PHE:CE1	6:F:1296:PHE:HE1	2.38	0.40
4:D:93:ASP:HB3	4:D:97:ARG:HH12	1.86	0.40
1:A:346:ARG:HD2	5:E:37:TYR:HB3	2.03	0.40
1:A:572:PHE:HB2	1:A:601:ILE:HD11	2.03	0.40
4:D:99:ASN:O	4:D:102:HIS:CE1	2.75	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	634/810 (78%)	599 (94%)	35 (6%)	0	100	100
2	B	365/392 (93%)	331 (91%)	34 (9%)	0	100	100
3	C	54/344 (16%)	53 (98%)	1 (2%)	0	100	100
4	D	211/245 (86%)	199 (94%)	11 (5%)	1 (0%)	24	52
5	E	81/123 (66%)	76 (94%)	4 (5%)	1 (1%)	10	33
6	F	224/353 (64%)	214 (96%)	10 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1569/2267 (69%)	1472 (94%)	95 (6%)	2 (0%)	49 76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	242	SER
5	E	108	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	542/689 (79%)	539 (99%)	3 (1%)	78 79
2	B	294/321 (92%)	292 (99%)	2 (1%)	76 77
3	C	41/276 (15%)	41 (100%)	0	100 100
4	D	178/204 (87%)	175 (98%)	3 (2%)	53 67
5	E	72/103 (70%)	72 (100%)	0	100 100
6	F	189/283 (67%)	189 (100%)	0	100 100
All	All	1316/1876 (70%)	1308 (99%)	8 (1%)	76 79

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	290	ILE
1	A	361	LYS
1	A	467	THR
2	B	63	LEU
2	B	144	GLN
4	D	31	PRO
4	D	130	VAL
4	D	136	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	GLN
1	A	179	GLN
1	A	181	ASN
1	A	217	GLN
1	A	244	GLN
1	A	384	GLN
1	A	411	GLN
1	A	445	GLN
1	A	466	GLN
1	A	475	ASN
1	A	506	ASN
1	A	534	ASN
1	A	538	ASN
1	A	542	GLN
1	A	555	HIS
1	A	561	GLN
1	A	563	ASN
1	A	693	GLN
1	A	775	GLN
2	B	58	ASN
2	B	165	HIS
2	B	225	GLN
2	B	230	GLN
2	B	234	GLN
2	B	266	ASN
2	B	287	ASN
2	B	302	ASN
2	B	324	HIS
2	B	334	ASN
2	B	347	HIS
2	B	359	GLN
4	D	50	GLN
4	D	60	ASN
4	D	99	ASN
4	D	102	HIS
4	D	158	GLN
5	E	52	GLN
5	E	83	HIS
5	E	89	GLN
6	F	1026	ASN
6	F	1071	HIS
6	F	1127	HIS
6	F	1128	HIS

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Mol	Chain	Res	Type
6	F	1160	HIS
6	F	1200	ASN
6	F	1230	GLN
6	F	1236	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

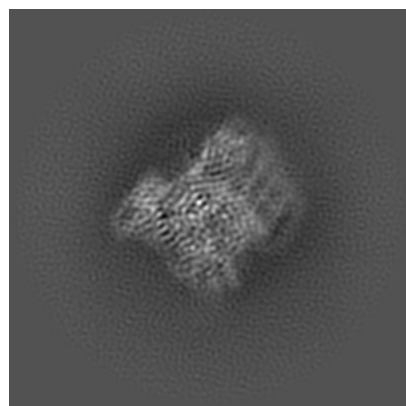
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-18507. These allow visual inspection of the internal detail of the map and identification of artifacts.

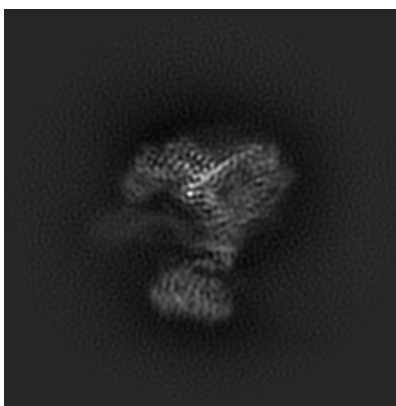
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

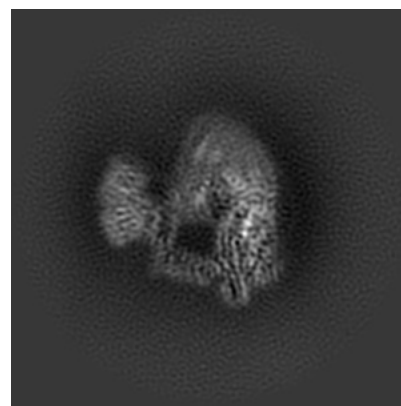
6.1.1 Primary map



X

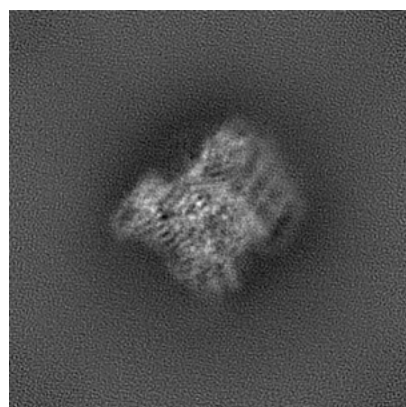


Y

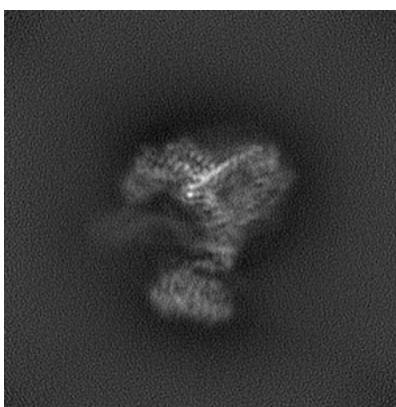


Z

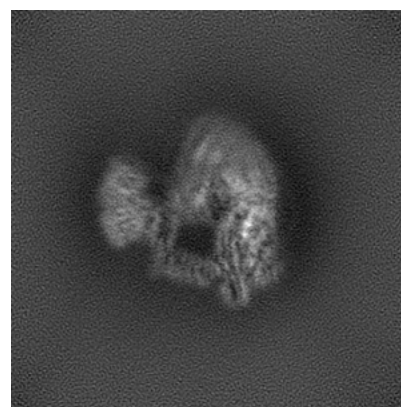
6.1.2 Raw map



X



Y

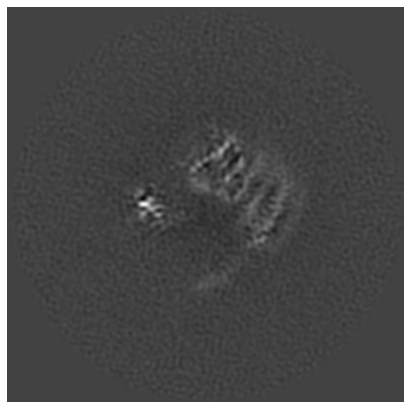


Z

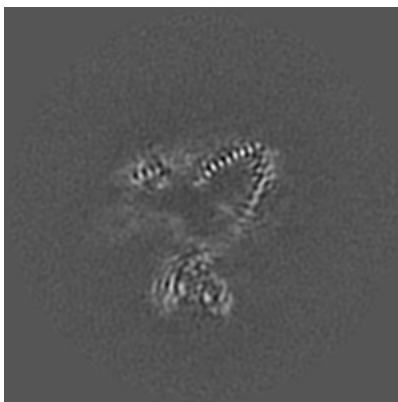
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

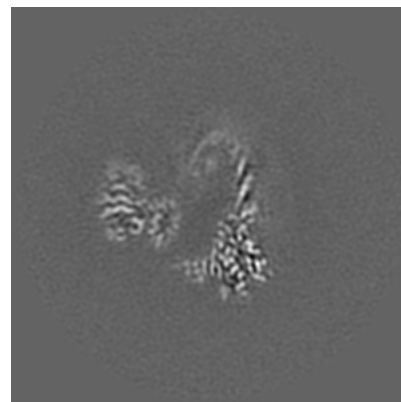
6.2.1 Primary map



X Index: 128



Y Index: 128

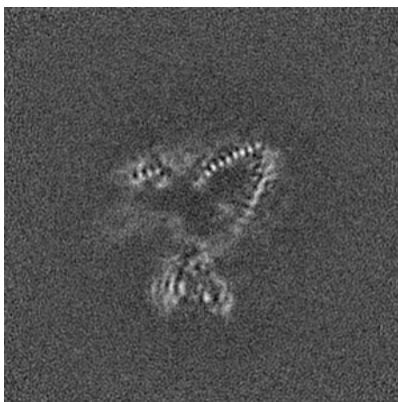


Z Index: 128

6.2.2 Raw map



X Index: 128



Y Index: 128

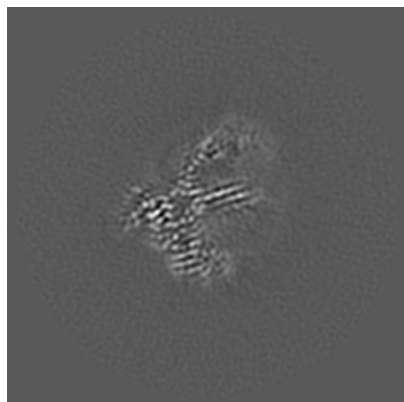


Z Index: 128

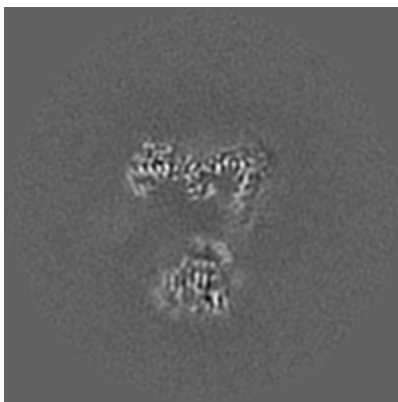
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

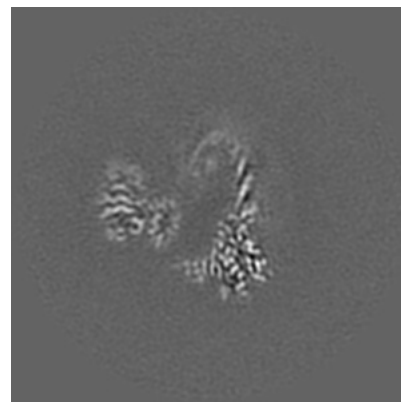
6.3.1 Primary map



X Index: 153

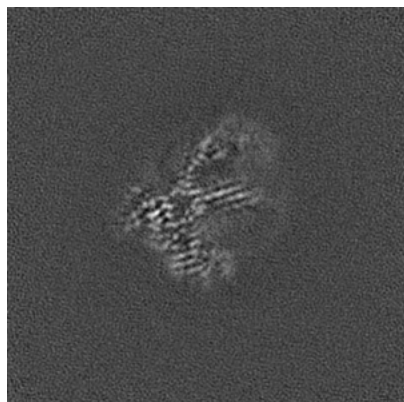


Y Index: 118

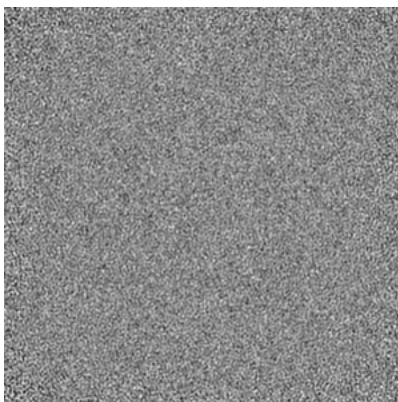


Z Index: 128

6.3.2 Raw map



X Index: 153



Y Index: 0

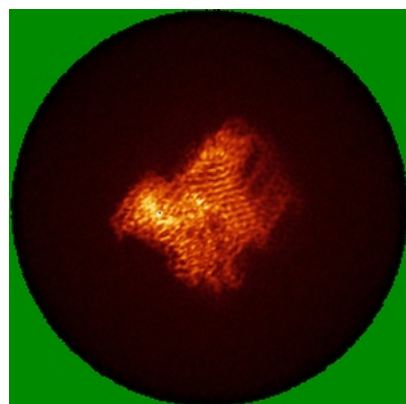


Z Index: 128

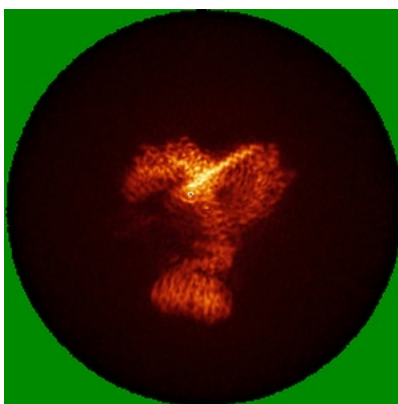
The images above show the largest variance slices of the map in three orthogonal directions.

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

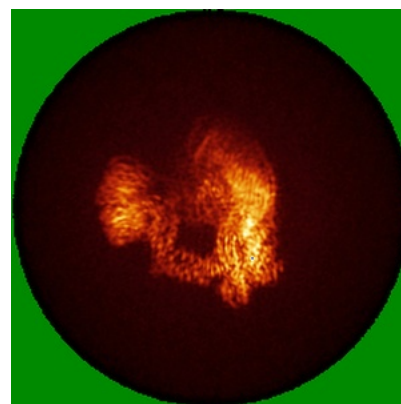
6.4.1 Primary map



X

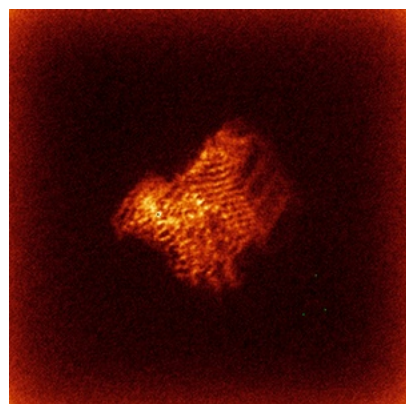


Y

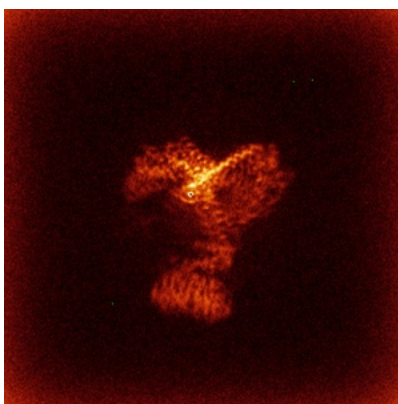


Z

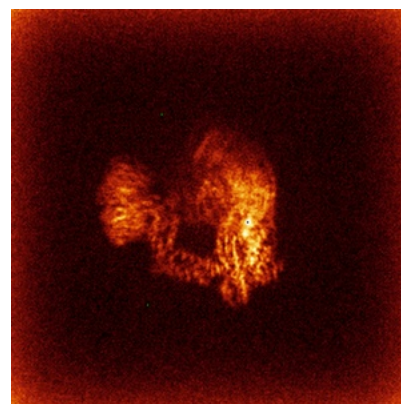
6.4.2 Raw map



X



Y

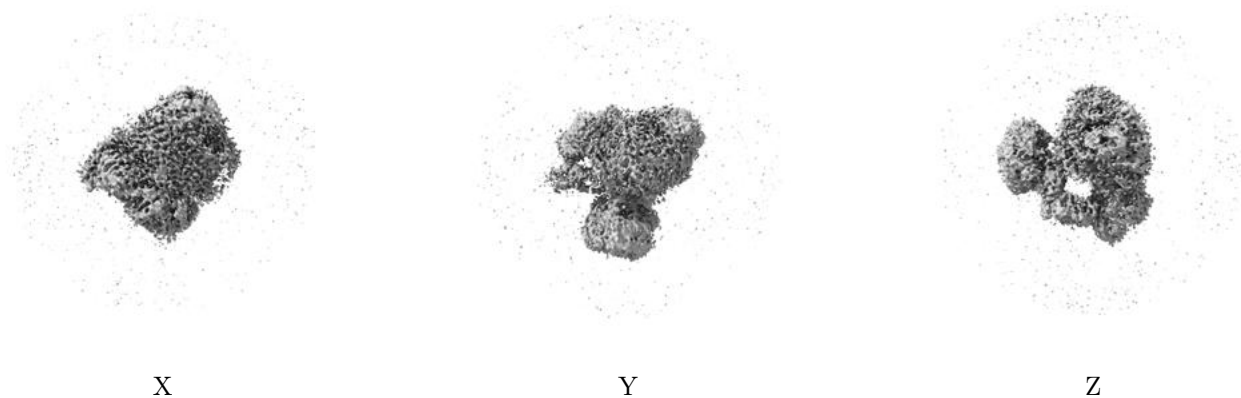


Z

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.

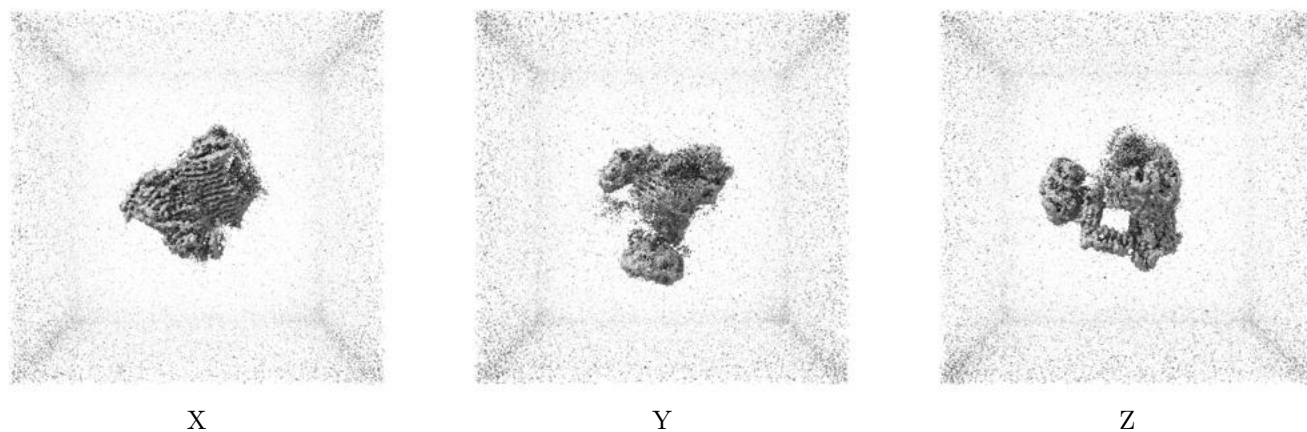
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. 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.

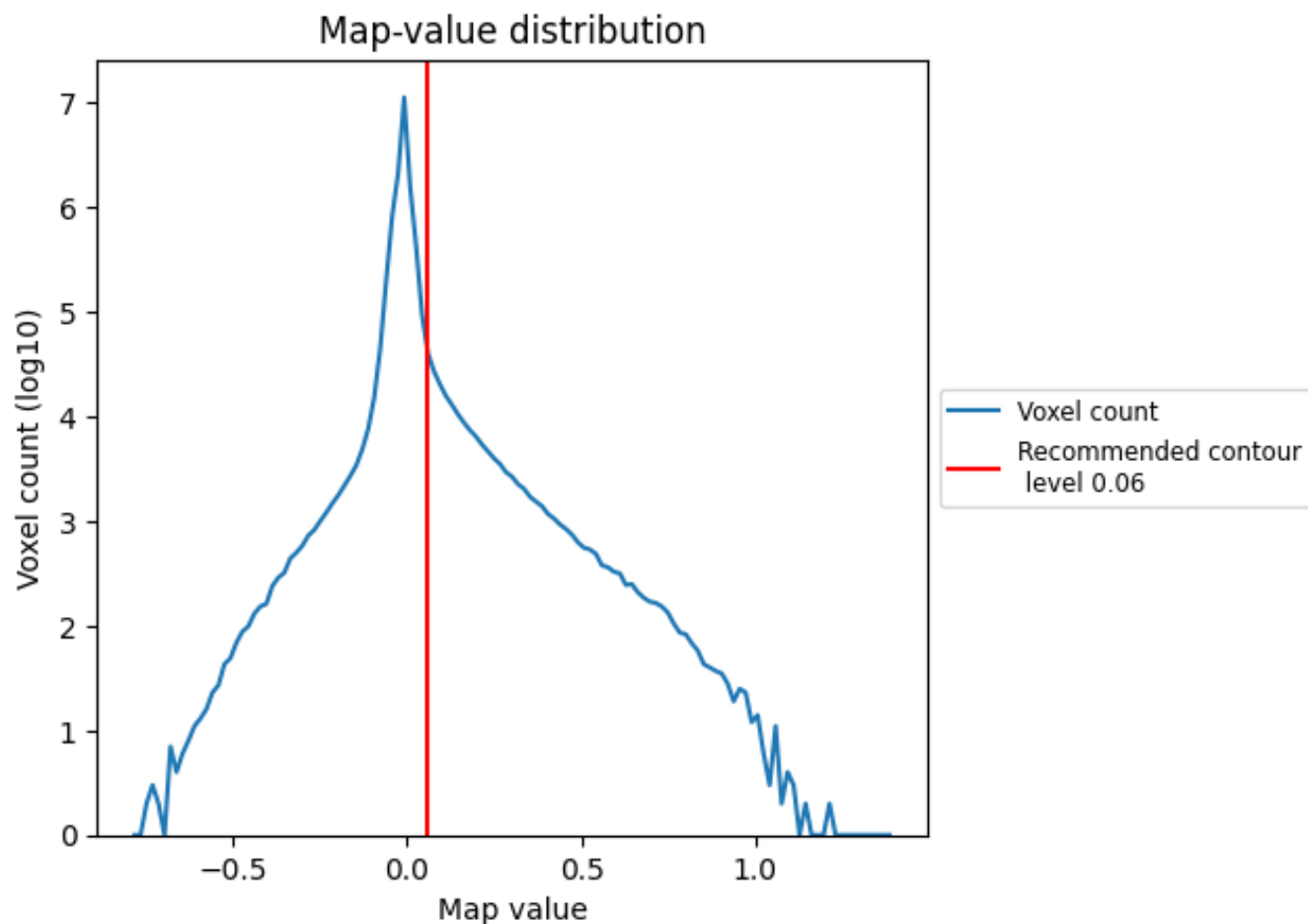
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

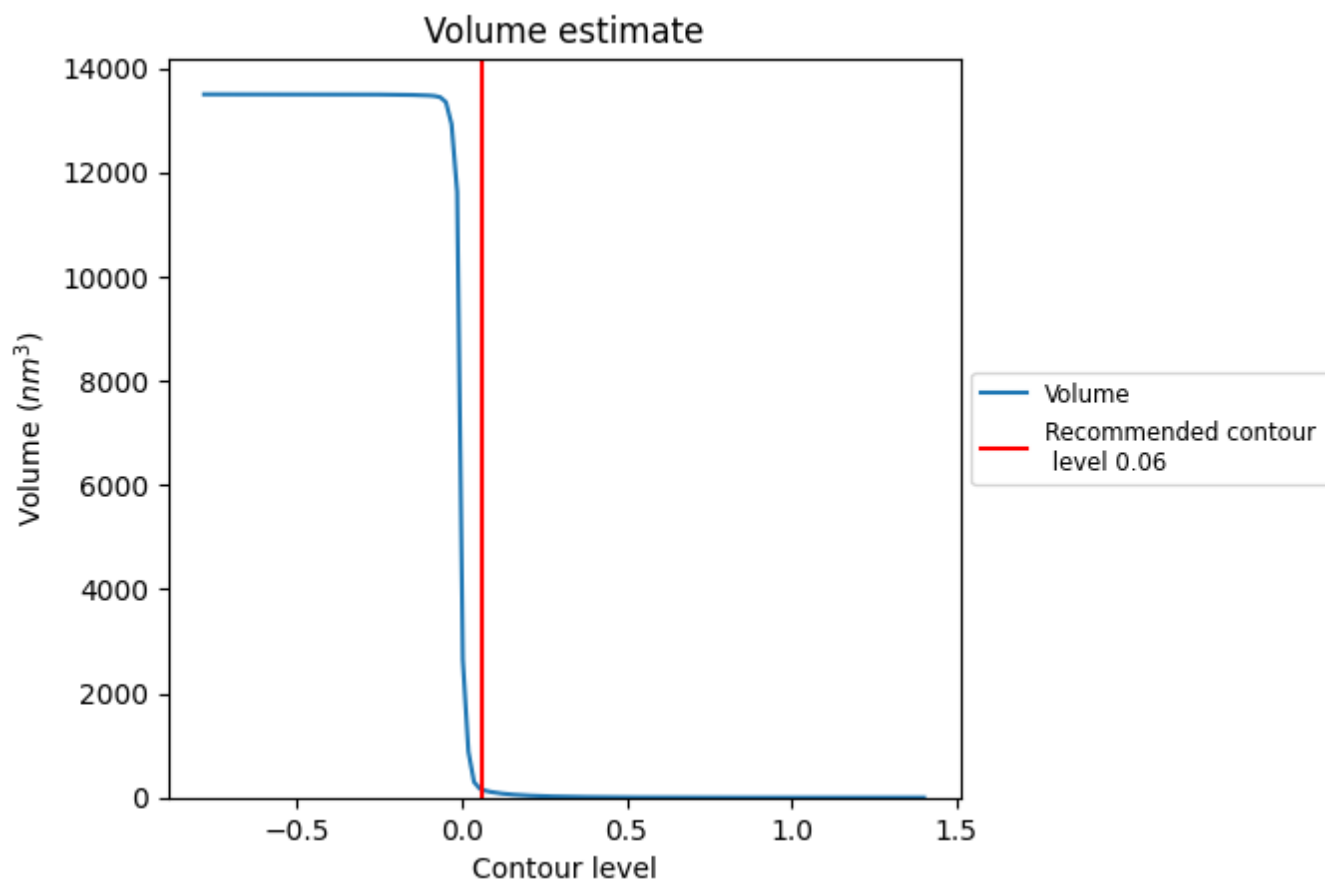
This section contains the results of statistical analysis of the map.

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.

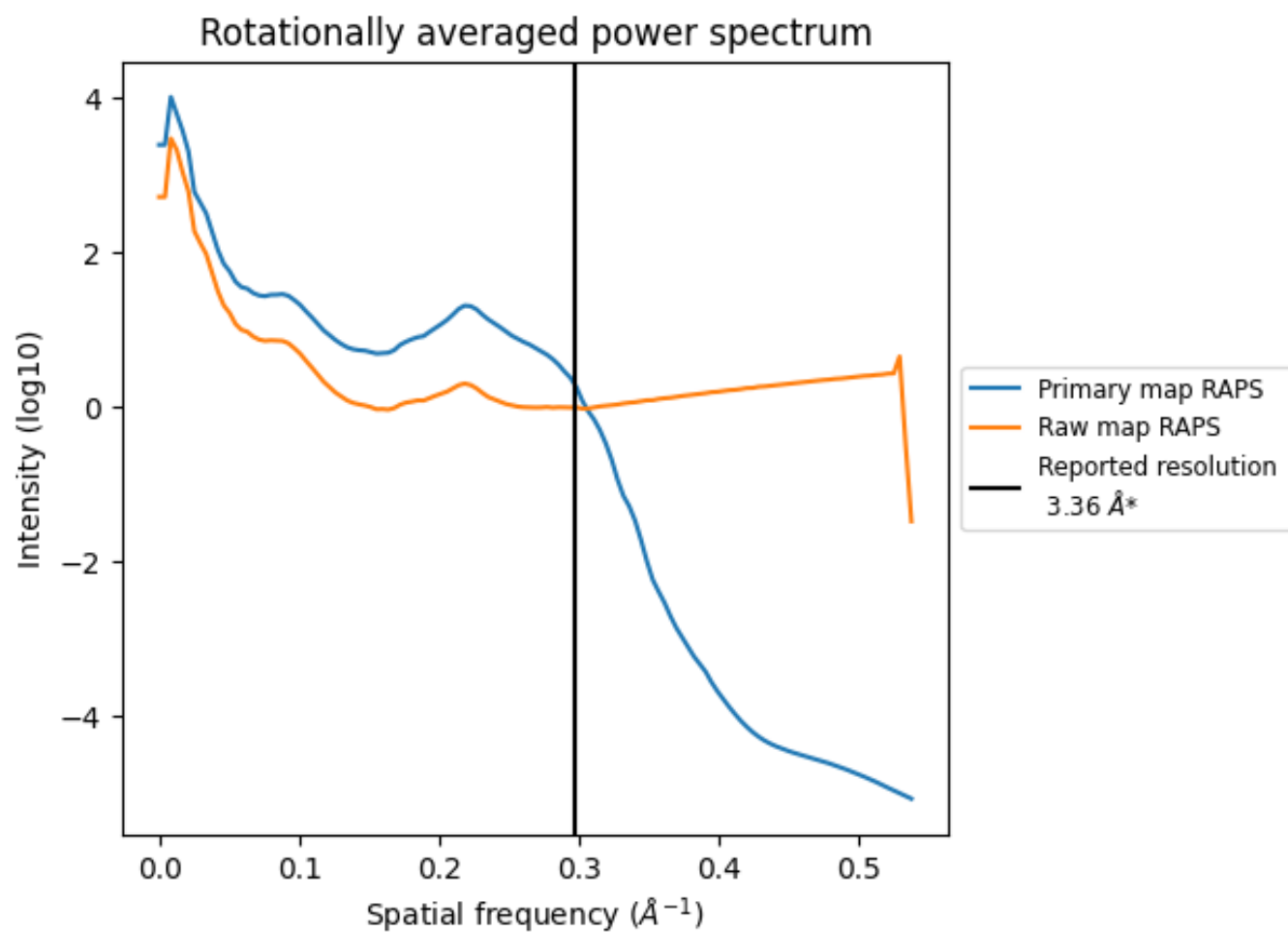
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 158 nm³; this corresponds to an approximate mass of 143 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 [i](#)

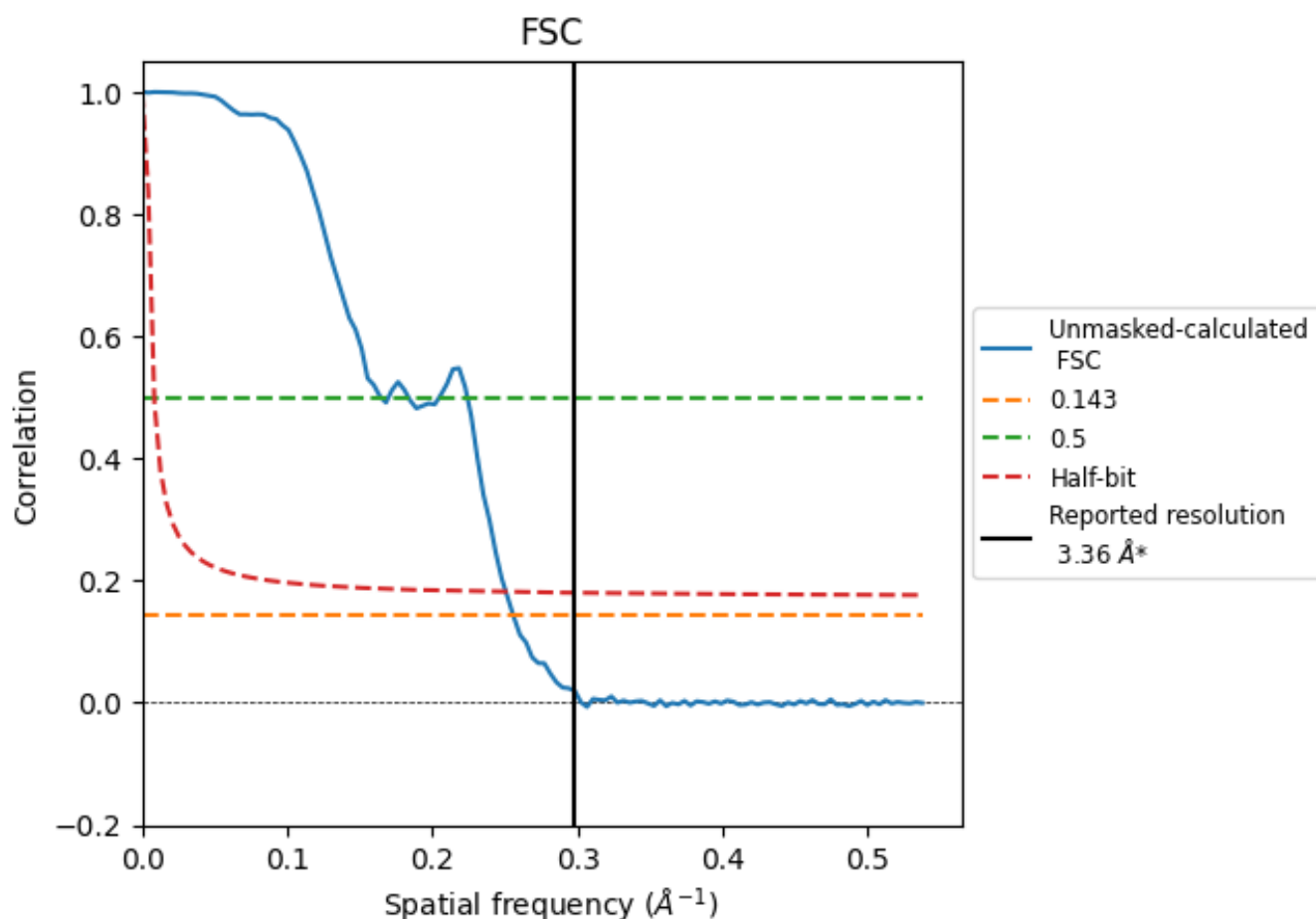


*Reported resolution corresponds to spatial frequency of 0.298 Å⁻¹

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 0.298 Å⁻¹

8.2 Resolution estimates [i](#)

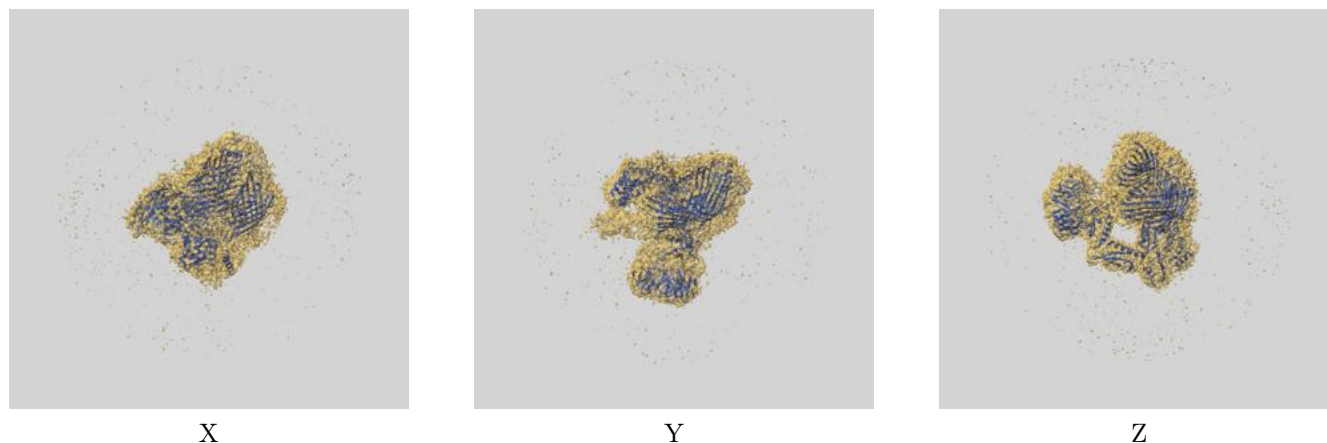
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.36	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.91	6.08	3.99

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.91 differs from the reported value 3.36 by more than 10 %

9 Map-model fit [i](#)

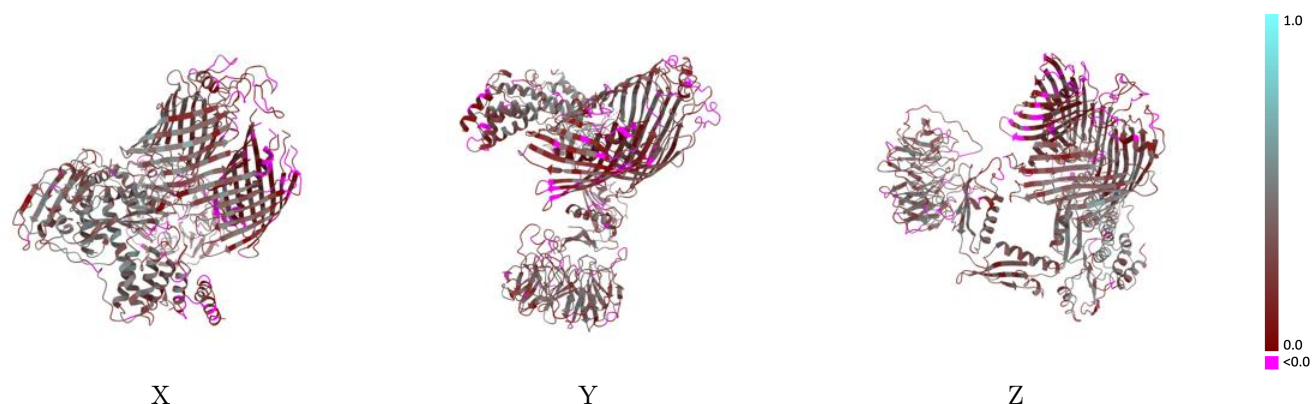
This section contains information regarding the fit between EMDB map EMD-18507 and PDB model 8QN4. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



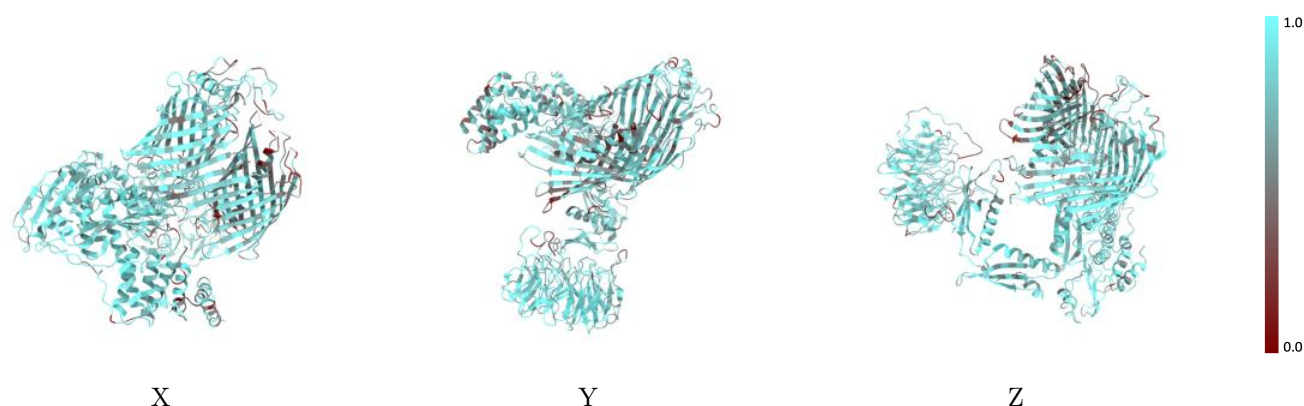
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



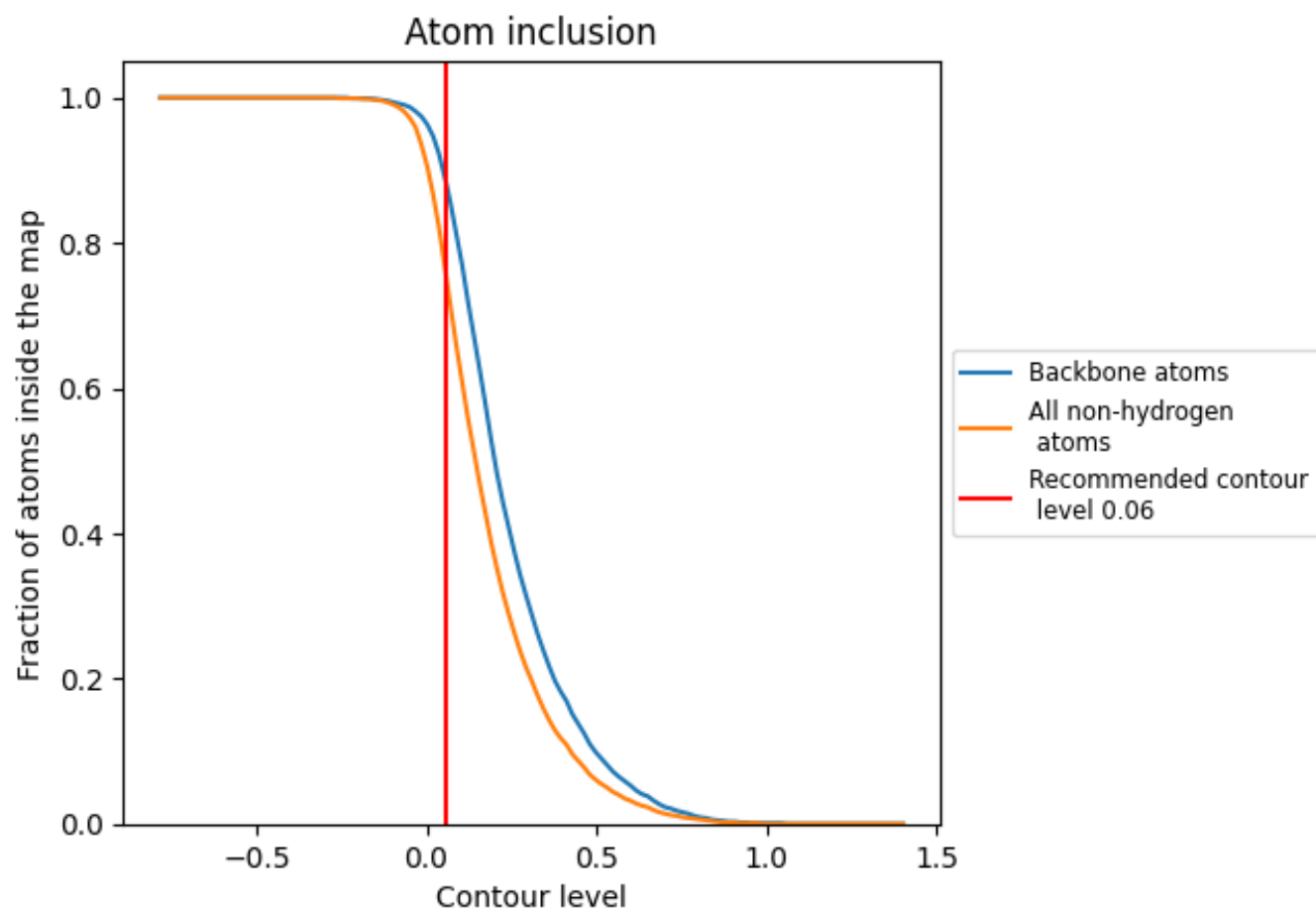
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 75% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7490	0.2710
A	0.7630	0.2920
B	0.7860	0.2770
C	0.7060	0.2560
D	0.7850	0.3110
E	0.8210	0.3390
F	0.6060	0.1450

1.0

0.0

<0.0