



Full wwPDB EM Validation Report ⓘ

Mar 7, 2026 – 06:16 AM UTC

PDB ID : 8TW9 / pdb_00008tw9
EMDB ID : EMD-41663
Title : Cryo-EM structure of S. cerevisiae PolE-Ctf18-8-1-DNA
Authors : Yuan, Z.; Georgescu, R.; O'Donnell, M.; Li, H.
Deposited on : 2023-08-20
Resolution : 3.60 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:

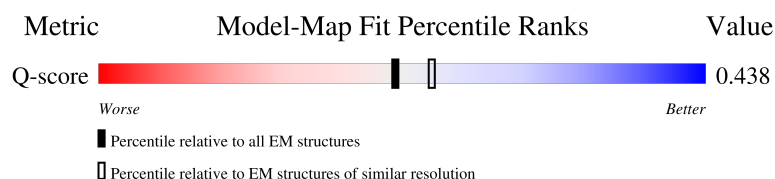
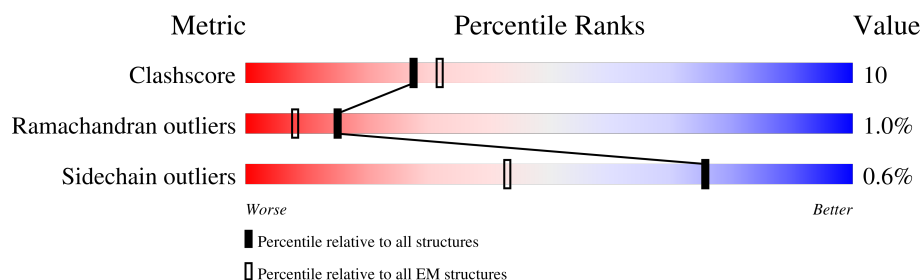
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
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 3.60 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	C	26	
2	P	9	
3	T	15	
4	E	2222	

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Mol	Chain	Length	Quality of chain
5	D	132	
6	B	380	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	SF4	E	2301	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 13586 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 Chromosome transmission fidelity protein 18.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	C	26	Total	C	N	O	0	0
			206	132	37	37		

- Molecule 2 is a DNA chain called Primer DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	P	9	Total	C	N	O	P	0	0
			183	88	29	57	9		

- Molecule 3 is a DNA chain called Template DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	T	15	Total	C	N	O	P	0	0
			303	147	54	88	14		

- Molecule 4 is a protein called DNA polymerase epsilon catalytic subunit A.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	1124	Total	C	N	O	S	0	0
			8919	5681	1507	1690	41		

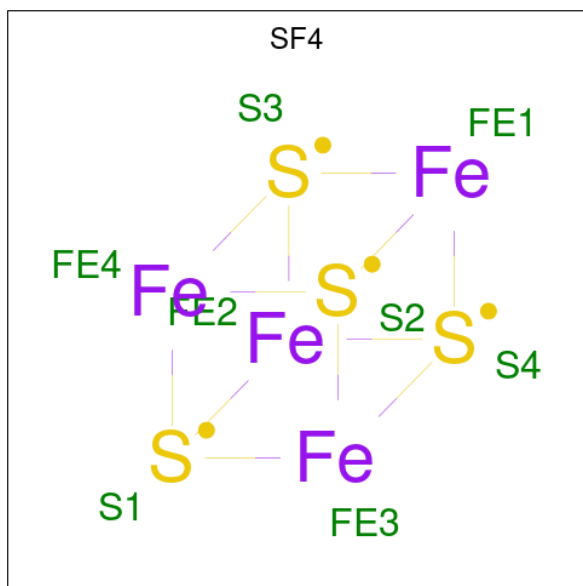
- Molecule 5 is a protein called Chromosome transmission fidelity protein 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	D	132	Total	C	N	O	S	0	0
			1014	642	171	194	7		

- Molecule 6 is a protein called Sister chromatid cohesion protein DCC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	B	374	Total	C	N	O	S	0	0
			2953	1893	489	557	14		

- Molecule 7 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).

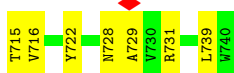


Mol	Chain	Residues	Atoms			AltConf
7	E	1	Total	Fe	S	0
			8	4	4	

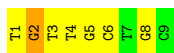
3 Residue-property plots

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.

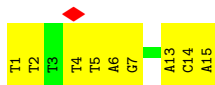
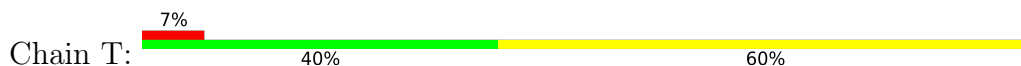
- Molecule 1: Chromosome transmission fidelity protein 18



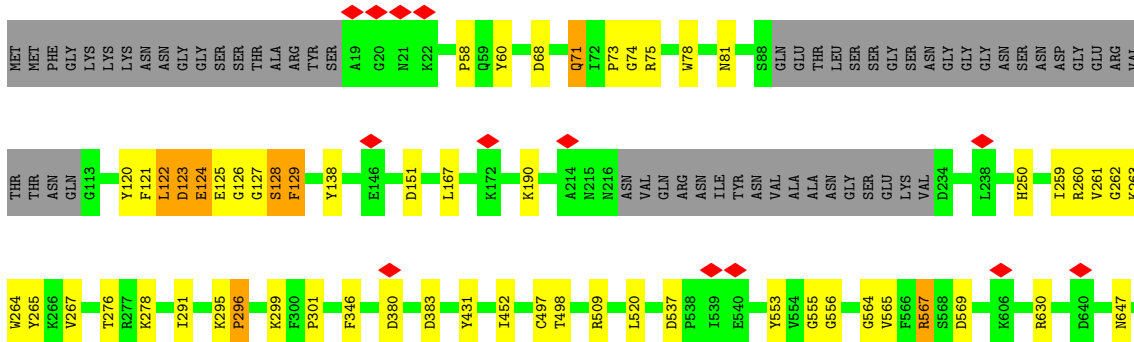
- Molecule 2: Primer DNA



- Molecule 3: Template DNA



- Molecule 4: DNA polymerase epsilon catalytic subunit A





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	385806	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.409	Depositor
Minimum map value	-0.591	Depositor
Average map value	0.009	Depositor
Map value standard deviation	0.060	Depositor
Recommended contour level	0.326	Depositor
Map size ($\text{\AA}$)	211.968, 211.968, 211.968	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.828, 0.828, 0.828	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SF4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	C	0.35	0/211	0.79	0/289
2	P	0.88	2/203 (1.0%)	0.73	0/311
3	T	0.73	0/339	0.66	0/521
4	E	0.82	1/9121 (0.0%)	1.29	29/12348 (0.2%)
5	D	0.33	0/1030	0.85	0/1387
6	B	0.56	3/3016 (0.1%)	0.93	13/4083 (0.3%)
All	All	0.74	6/13920 (0.0%)	1.16	42/18939 (0.2%)

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
4	E	0	6
6	B	0	2
All	All	0	8

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	354	ARG	C-N	14.18	1.54	1.33
4	E	766	GLU	C-N	10.87	1.56	1.33
6	B	297	ASP	C-O	7.92	1.32	1.23
6	B	355	GLY	C-N	-7.54	1.21	1.33
2	P	2	DG	C1'-N9	-5.81	1.34	1.46
2	P	8	DG	C1'-N9	-5.62	1.35	1.46

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	766	GLU	O-C-N	-9.10	111.66	122.95
4	E	1060	GLU	N-CA-C	-8.71	100.61	111.75
4	E	122	LEU	N-CA-C	-8.60	100.75	111.75
6	B	355	GLY	CA-C-N	-8.13	108.23	123.27
6	B	355	GLY	C-N-CA	-8.13	108.23	123.27
4	E	656	PRO	N-CA-C	-7.68	103.77	114.98
6	B	352	ASN	O-C-N	7.62	133.05	123.14
4	E	151	ASP	CA-CB-CG	7.57	120.17	112.60
4	E	1058	GLU	N-CA-C	-7.27	103.35	111.71
4	E	877	ASP	CA-CB-CG	6.83	119.43	112.60
4	E	120	TYR	CA-C-N	-6.52	113.86	123.11
4	E	120	TYR	C-N-CA	-6.52	113.86	123.11
4	E	1075	ASP	CA-CB-CG	6.47	119.07	112.60
6	B	303	GLY	N-CA-C	-6.47	105.70	113.99
4	E	1059	TYR	CA-CB-CG	-6.41	102.37	113.90
4	E	380	ASP	CA-CB-CG	6.15	118.75	112.60
4	E	835	MET	N-CA-C	-5.90	104.28	112.30
4	E	671	ASN	OD1-CG-ND2	-5.81	116.79	122.60
4	E	124	GLU	N-CA-C	-5.76	101.99	110.52
4	E	296	PRO	N-CA-CB	5.70	106.18	103.22
4	E	830	PHE	N-CA-C	-5.63	102.08	111.37
6	B	87	LEU	CA-CB-CG	5.62	135.98	116.30
4	E	71	GLN	N-CA-C	-5.53	105.66	112.90
4	E	129	PHE	CA-C-N	-5.47	112.78	122.74
4	E	129	PHE	C-N-CA	-5.47	112.78	122.74
4	E	834	VAL	CA-C-O	-5.47	113.94	120.78
6	B	375	THR	CA-C-N	-5.43	116.34	122.48
6	B	375	THR	C-N-CA	-5.43	116.34	122.48
6	B	185	LEU	CA-CB-CG	5.40	135.20	116.30
4	E	1076	PHE	CA-CB-CG	5.21	119.01	113.80
6	B	352	ASN	CA-C-N	-5.19	111.62	121.54
6	B	352	ASN	C-N-CA	-5.19	111.62	121.54
4	E	817	ASP	O-C-N	5.19	127.62	122.12
6	B	351	LEU	CA-C-N	-5.17	114.94	122.44
6	B	351	LEU	C-N-CA	-5.17	114.94	122.44
4	E	250	HIS	CB-CG-CD2	-5.14	124.52	131.20
4	E	1105	ARG	NE-CZ-NH2	5.14	123.83	119.20
6	B	362	ILE	CA-C-O	-5.14	116.06	121.41
4	E	537	ASP	CA-CB-CG	5.14	117.74	112.60
4	E	766	GLU	CA-C-N	5.04	134.09	121.80
4	E	766	GLU	C-N-CA	5.04	134.09	121.80
4	E	968	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
6	B	305	HIS	Mainchain
6	B	355	GLY	Mainchain
4	E	123	ASP	Mainchain
4	E	509	ARG	Sidechain
4	E	567	ARG	Sidechain
4	E	630	ARG	Sidechain
4	E	766	GLU	Mainchain
4	E	81	ASN	Mainchain

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	206	0	183	8	0
2	P	183	0	102	6	0
3	T	303	0	172	15	0
4	E	8919	0	8572	159	0
5	D	1014	0	1007	30	0
6	B	2953	0	2913	80	0
7	E	8	0	0	6	0
All	All	13586	0	12949	257	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 10.

All (257) 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)
4:E:656:PRO:HB3	4:E:843:SER:HB2	1.46	0.95
3:T:5:DT:H4'	4:E:835:MET:CB	2.08	0.83
4:E:60:TYR:CB	4:E:73:PRO:HG3	2.13	0.79
2:P:2:DG:H2'	2:P:3:DT:H71	1.68	0.76
4:E:671:ASN:HB2	7:E:2301:SF4:S3	2.26	0.75
4:E:845:GLU:O	4:E:847:ALA:N	2.21	0.73
1:C:722:TYR:H	4:E:1041:ASP:CB	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:784:PHE:CD1	4:E:816:TYR:HD1	2.07	0.73
4:E:78:TRP:CZ2	4:E:262:GLY:HA2	2.24	0.72
3:T:4:DT:O2	4:E:832:GLY:HA3	1.90	0.72
6:B:335:LEU:HB2	6:B:339:TRP:HZ2	1.56	0.71
4:E:827:LEU:HA	4:E:830:PHE:HD2	1.55	0.70
4:E:784:PHE:HD1	4:E:816:TYR:CD1	2.09	0.70
6:B:303:GLY:O	6:B:346:PRO:HG3	1.91	0.70
4:E:665:CYS:HA	7:E:2301:SF4:S2	2.30	0.70
4:E:784:PHE:CD1	4:E:816:TYR:CD1	2.79	0.69
4:E:78:TRP:HB2	4:E:264:TRP:CD2	2.29	0.68
4:E:567:ARG:HH11	4:E:567:ARG:HG2	1.58	0.68
5:D:103:ILE:HD11	6:B:15:TYR:HB3	1.75	0.68
4:E:125:GLU:C	6:B:368:ARG:O	2.37	0.68
4:E:259:ILE:HA	4:E:265:TYR:CZ	2.30	0.66
4:E:78:TRP:HE1	4:E:262:GLY:C	2.05	0.65
4:E:124:GLU:O	4:E:126:GLY:N	2.30	0.65
4:E:260:ARG:N	4:E:265:TYR:OH	2.29	0.65
4:E:845:GLU:O	4:E:848:GLY:N	2.29	0.65
4:E:986:LEU:O	4:E:988:ARG:N	2.30	0.64
4:E:677:CYS:O	4:E:764:GLN:HG2	1.98	0.64
4:E:986:LEU:C	4:E:988:ARG:H	2.06	0.64
4:E:781:ARG:HE	4:E:824:LYS:HB2	1.61	0.64
3:T:1:DT:H4'	3:T:2:DT:OP1	1.97	0.64
4:E:658:SER:O	4:E:659:ILE:C	2.41	0.64
4:E:121:PHE:C	4:E:123:ASP:N	2.50	0.63
4:E:295:LYS:HB3	4:E:301:PRO:HG3	1.80	0.63
4:E:1155:GLN:HA	4:E:1159:THR:OG1	1.98	0.63
4:E:1059:TYR:C	4:E:1061:GLY:N	2.54	0.62
6:B:302:ARG:CB	6:B:307:LYS:HE3	2.29	0.62
5:D:15:THR:HG21	6:B:29:GLN:HE22	1.64	0.62
6:B:332:LEU:O	6:B:339:TRP:NE1	2.32	0.62
3:T:7:DG:H21	4:E:967:LYS:HD2	1.65	0.62
4:E:656:PRO:HD3	4:E:770:TYR:CE2	2.34	0.61
6:B:341:LEU:C	6:B:343:ASP:N	2.58	0.60
4:E:829:SER:O	4:E:830:PHE:C	2.44	0.60
2:P:5:DG:H2''	2:P:6:DC:H5'	1.84	0.60
4:E:657:ASP:O	4:E:766:GLU:HG2	2.02	0.59
4:E:1059:TYR:O	4:E:1060:GLU:C	2.44	0.59
6:B:335:LEU:HB2	6:B:339:TRP:CZ2	2.37	0.59
2:P:1:DT:OP1	4:E:1179:TRP:HZ2	1.86	0.59
5:D:12:GLN:NE2	6:B:28:ILE:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:7:DG:OP2	4:E:686:ARG:NH1	2.33	0.59
6:B:305:HIS:CB	6:B:314:GLN:O	2.51	0.58
6:B:363:MET:HA	6:B:366:ALA:O	2.03	0.58
1:C:731:ARG:HH21	6:B:61:LYS:HG3	1.69	0.58
6:B:351:LEU:O	6:B:352:ASN:OD1	2.21	0.58
1:C:715:THR:HG23	1:C:716:VAL:HG23	1.86	0.57
4:E:827:LEU:HA	4:E:830:PHE:CD2	2.38	0.57
4:E:262:GLY:C	4:E:498:THR:HG22	2.29	0.57
4:E:799:ASP:OD1	4:E:799:ASP:N	2.38	0.57
6:B:331:VAL:HA	6:B:334:ARG:HD3	1.86	0.57
6:B:262:GLY:HA2	6:B:284:TRP:HE1	1.68	0.57
3:T:5:DT:C4'	4:E:835:MET:CB	2.82	0.56
6:B:341:LEU:C	6:B:343:ASP:H	2.12	0.56
6:B:40:LYS:NZ	6:B:152:GLU:O	2.40	0.55
4:E:807:ASP:O	4:E:811:LYS:HG3	2.06	0.55
4:E:832:GLY:O	4:E:833:TYR:C	2.50	0.55
4:E:262:GLY:O	4:E:498:THR:HG22	2.07	0.55
6:B:358:ILE:O	6:B:362:ILE:HG13	2.07	0.55
4:E:980:GLU:O	4:E:981:LEU:HD23	2.07	0.54
6:B:341:LEU:O	6:B:344:ILE:N	2.34	0.54
6:B:341:LEU:O	6:B:343:ASP:N	2.39	0.54
4:E:125:GLU:CB	6:B:369:LYS:HA	2.38	0.54
4:E:1025:ARG:CD	5:D:89:ARG:NH2	2.71	0.54
4:E:972:PHE:CZ	4:E:1003:PHE:O	2.61	0.54
3:T:7:DG:H4'	4:E:556:GLY:O	2.08	0.53
4:E:260:ARG:HB2	4:E:263:LYS:HD2	1.90	0.53
5:D:89:ARG:N	5:D:130:LEU:O	2.41	0.53
1:C:728:ASN:HD21	6:B:63:ARG:HD3	1.74	0.53
4:E:829:SER:O	4:E:832:GLY:N	2.41	0.53
6:B:297:ASP:C	6:B:299:ASP:N	2.66	0.53
4:E:121:PHE:O	4:E:122:LEU:C	2.51	0.53
6:B:330:LYS:O	6:B:334:ARG:HG3	2.09	0.53
4:E:1058:GLU:C	4:E:1059:TYR:HD2	2.17	0.52
6:B:27:ILE:HG23	6:B:33:GLN:HE22	1.73	0.52
4:E:656:PRO:HB3	4:E:843:SER:CB	2.30	0.52
4:E:1019:VAL:CG1	4:E:1158:ILE:HG12	2.40	0.52
6:B:197:MET:HE1	6:B:292:PHE:HA	1.91	0.52
4:E:802:ASP:HA	4:E:806:ARG:HE	1.74	0.52
6:B:303:GLY:O	6:B:346:PRO:CG	2.56	0.52
4:E:78:TRP:HZ2	4:E:262:GLY:HA2	1.73	0.51
4:E:296:PRO:HB2	4:E:299:LYS:HD2	1.90	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:658:SER:O	4:E:660:LYS:N	2.43	0.51
4:E:791:TRP:HB3	4:E:813:ILE:HG12	1.93	0.51
1:C:729:ALA:HA	4:E:452:ILE:HD12	1.91	0.51
4:E:1025:ARG:HD3	5:D:89:ARG:NH2	2.25	0.51
4:E:1044:SER:O	4:E:1047:CYS:HB3	2.10	0.51
4:E:291:ILE:HG12	4:E:383:ASP:OD1	2.11	0.51
4:E:969:TYR:O	4:E:982:LYS:N	2.44	0.51
6:B:309:THR:HG23	6:B:311:LYS:H	1.75	0.51
4:E:830:PHE:O	4:E:834:VAL:HG13	2.11	0.51
4:E:1059:TYR:N	4:E:1059:TYR:CD2	2.79	0.51
5:D:23:THR:OG1	5:D:24:VAL:N	2.44	0.50
4:E:276:THR:C	4:E:278:LYS:H	2.19	0.50
4:E:647:ASN:HA	4:E:650:THR:HG22	1.93	0.50
4:E:1055:THR:N	4:E:1084:ASP:O	2.44	0.50
5:D:106:PHE:HE2	6:B:16:LYS:HD2	1.76	0.50
6:B:297:ASP:C	6:B:299:ASP:H	2.18	0.50
4:E:969:TYR:H	4:E:982:LYS:CB	2.24	0.50
5:D:106:PHE:N	6:B:14:SER:OG	2.44	0.50
6:B:359:ASP:O	6:B:363:MET:HG2	2.11	0.50
4:E:123:ASP:CB	4:E:129:PHE:CE2	2.95	0.50
3:T:6:DA:O3'	4:E:555:GLY:N	2.38	0.50
4:E:78:TRP:CD1	4:E:264:TRP:CD1	2.99	0.50
1:C:729:ALA:O	6:B:63:ARG:NH1	2.45	0.49
6:B:306:PHE:HB2	6:B:339:TRP:CH2	2.47	0.49
4:E:124:GLU:HB3	4:E:264:TRP:HZ3	1.76	0.49
4:E:1057:LYS:C	4:E:1059:TYR:N	2.67	0.49
4:E:845:GLU:O	4:E:846:MET:C	2.55	0.49
5:D:2:PRO:HG3	6:B:44:LYS:HG2	1.94	0.49
4:E:78:TRP:HD1	4:E:264:TRP:CD1	2.31	0.49
4:E:263:LYS:HE3	4:E:498:THR:O	2.13	0.49
4:E:1046:ILE:O	4:E:1047:CYS:C	2.55	0.49
6:B:239:ALA:HB2	6:B:252:LEU:HD23	1.95	0.49
6:B:261:TYR:OH	6:B:284:TRP:O	2.22	0.49
6:B:363:MET:O	6:B:364:LYS:C	2.55	0.49
5:D:71:GLY:CA	5:D:83:PHE:O	2.61	0.48
6:B:352:ASN:C	6:B:354:ARG:H	2.21	0.48
3:T:13:DA:H2''	3:T:14:DC:C6	2.49	0.48
4:E:986:LEU:C	4:E:988:ARG:N	2.66	0.48
4:E:78:TRP:NE1	4:E:262:GLY:C	2.71	0.48
4:E:832:GLY:O	4:E:834:VAL:N	2.47	0.48
5:D:13:LYS:O	5:D:16:GLN:NE2	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:3:SER:HB2	6:B:38:ARG:HE	1.78	0.47
4:E:770:TYR:OH	4:E:834:VAL:HG12	2.13	0.47
2:P:3:DT:H2''	2:P:4:DT:O5'	2.15	0.47
4:E:1158:ILE:O	4:E:1159:THR:C	2.58	0.47
5:D:73:LEU:O	6:B:2:SER:OG	2.28	0.47
6:B:18:ILE:HG22	6:B:108:PHE:HB2	1.95	0.47
3:T:4:DT:O2	4:E:832:GLY:CA	2.62	0.47
6:B:302:ARG:C	6:B:304:TYR:N	2.71	0.47
4:E:845:GLU:C	4:E:847:ALA:N	2.72	0.47
4:E:1056:LEU:HD22	4:E:1084:ASP:O	2.15	0.47
4:E:777:PHE:HA	4:E:780:ARG:HG2	1.96	0.46
4:E:78:TRP:HA	4:E:264:TRP:HA	1.97	0.46
4:E:1157:ILE:O	4:E:1161:PRO:HG2	2.15	0.46
4:E:815:LEU:O	4:E:818:SER:OG	2.28	0.46
4:E:840:ARG:C	4:E:842:TYR:N	2.72	0.46
6:B:298:ILE:HA	6:B:301:LEU:CB	2.45	0.46
1:C:739:LEU:HD22	5:D:106:PHE:CE2	2.50	0.46
4:E:129:PHE:O	4:E:129:PHE:CD1	2.69	0.46
4:E:957:SER:HB2	4:E:1164:LEU:CB	2.46	0.46
5:D:11:TRP:NE1	5:D:117:ASP:OD1	2.42	0.46
4:E:1059:TYR:HD2	4:E:1059:TYR:N	2.13	0.46
4:E:78:TRP:HB2	4:E:264:TRP:CE3	2.51	0.46
5:D:73:LEU:HD23	6:B:5:LEU:HD21	1.98	0.46
4:E:968:ARG:HA	4:E:982:LYS:CB	2.46	0.46
6:B:340:ASP:HB2	6:B:343:ASP:HB2	1.98	0.46
3:T:6:DA:H5'	4:E:553:TYR:CE2	2.51	0.46
6:B:186:PHE:CG	6:B:264:GLN:HG3	2.51	0.46
6:B:212:HIS:HB2	6:B:231:ILE:HG13	1.98	0.46
4:E:792:LYS:HA	4:E:795:LEU:HD23	1.98	0.45
4:E:569:ASP:O	4:E:1010:ASP:O	2.34	0.45
4:E:784:PHE:CE1	4:E:816:TYR:CD1	3.04	0.45
4:E:1036:MET:CB	5:D:87:LYS:O	2.64	0.45
4:E:668:CYS:HB3	7:E:2301:SF4:S3	2.56	0.45
4:E:671:ASN:N	7:E:2301:SF4:S3	2.90	0.45
4:E:763:CYS:SG	7:E:2301:SF4:S1	3.14	0.45
6:B:242:GLU:N	6:B:249:THR:O	2.46	0.45
3:T:5:DT:H5'	4:E:835:MET:O	2.16	0.45
4:E:678:ALA:HB2	7:E:2301:SF4:S4	2.57	0.45
4:E:564:GLY:HA2	4:E:1161:PRO:C	2.42	0.45
3:T:7:DG:P	4:E:555:GLY:H	2.39	0.45
4:E:78:TRP:HB2	4:E:264:TRP:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:1025:ARG:HD2	5:D:89:ARG:NH2	2.31	0.45
6:B:362:ILE:O	6:B:366:ALA:O	2.35	0.45
4:E:520:LEU:CD1	4:E:833:TYR:CE2	3.00	0.45
5:D:73:LEU:HB3	6:B:5:LEU:HD11	1.98	0.45
5:D:133:MET:HA	6:B:89:PHE:HE2	1.81	0.45
4:E:431:TYR:CG	4:E:431:TYR:O	2.71	0.44
4:E:138:TYR:CG	4:E:190:LYS:HE3	2.52	0.44
3:T:6:DA:H5'	4:E:553:TYR:CZ	2.53	0.44
5:D:94:VAL:HG22	5:D:123:VAL:HG12	2.00	0.44
4:E:970:ALA:HB2	4:E:984:PHE:CZ	2.52	0.44
2:P:2:DG:H2'	2:P:3:DT:C6	2.53	0.44
2:P:5:DG:H2''	2:P:6:DC:C5'	2.47	0.44
4:E:123:ASP:CB	4:E:129:PHE:HE2	2.30	0.44
4:E:1064:SER:O	4:E:1068:THR:N	2.49	0.44
6:B:72:MET:HE2	6:B:72:MET:HB2	1.82	0.44
6:B:218:ASP:OD1	6:B:218:ASP:N	2.44	0.44
4:E:956:PRO:HB3	4:E:1156:LYS:O	2.18	0.43
3:T:14:DC:H2''	3:T:15:DA:C8	2.53	0.43
4:E:812:MET:HG3	4:E:815:LEU:HD23	2.00	0.43
5:D:89:ARG:HB2	5:D:132:ILE:HA	2.00	0.43
4:E:267:VAL:HG13	4:E:267:VAL:O	2.18	0.43
4:E:1052:MET:HB3	4:E:1059:TYR:CZ	2.53	0.43
4:E:1170:PRO:C	4:E:1172:PRO:HD3	2.42	0.43
6:B:341:LEU:O	6:B:342:GLU:C	2.62	0.43
5:D:31:MET:HE2	5:D:123:VAL:HG13	2.00	0.43
4:E:843:SER:O	4:E:844:MET:C	2.62	0.43
5:D:73:LEU:HD11	6:B:3:ILE:HD12	2.01	0.43
4:E:58:PRO:HD2	4:E:74:GLY:HA2	2.01	0.43
4:E:1160:ILE:O	4:E:1161:PRO:C	2.61	0.43
6:B:28:ILE:HD13	6:B:28:ILE:HA	1.82	0.43
4:E:833:TYR:O	4:E:836:ARG:N	2.39	0.43
4:E:564:GLY:HA2	4:E:1161:PRO:O	2.19	0.43
4:E:810:LYS:HA	4:E:813:ILE:HD12	2.00	0.43
4:E:565:VAL:O	4:E:565:VAL:HG13	2.19	0.42
6:B:220:GLU:HG3	6:B:223:PHE:HB2	2.00	0.42
4:E:993:GLN:HB3	4:E:1045:LEU:O	2.19	0.42
6:B:318:LYS:HE3	6:B:318:LYS:HB2	1.87	0.42
4:E:824:LYS:HG2	4:E:828:ASN:HD21	1.85	0.42
4:E:1052:MET:HE1	4:E:1055:THR:HG23	2.00	0.42
6:B:361:PHE:O	6:B:364:LYS:HB3	2.19	0.42
4:E:812:MET:HA	4:E:815:LEU:HB3	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:1036:MET:CB	5:D:131:PRO:HG3	2.50	0.42
4:E:276:THR:C	4:E:278:LYS:N	2.75	0.42
6:B:51:LEU:HB2	6:B:60:LEU:HD11	2.02	0.42
6:B:336:GLN:O	6:B:339:TRP:CD1	2.72	0.42
6:B:356:MET:HE2	6:B:360:SER:OG	2.20	0.42
4:E:127:GLY:O	4:E:128:SER:CB	2.67	0.42
6:B:362:ILE:HB	6:B:363:MET:HE2	2.00	0.42
4:E:75:ARG:CZ	4:E:129:PHE:CD2	3.03	0.42
4:E:794:ASN:HA	4:E:797:LYS:HG2	2.01	0.42
4:E:834:VAL:O	4:E:834:VAL:HG23	2.20	0.41
4:E:125:GLU:CB	6:B:368:ARG:O	2.67	0.41
5:D:82:LEU:HD23	5:D:90:LEU:HD12	2.02	0.41
6:B:74:GLU:HA	6:B:98:VAL:HG12	2.01	0.41
6:B:180:LEU:HD13	6:B:180:LEU:HA	1.74	0.41
6:B:254:ILE:HD13	6:B:257:ILE:HD12	2.02	0.41
4:E:833:TYR:O	4:E:835:MET:N	2.53	0.41
6:B:331:VAL:O	6:B:335:LEU:HG	2.20	0.41
1:C:729:ALA:HA	4:E:452:ILE:CD1	2.50	0.41
5:D:98:ASP:N	5:D:98:ASP:OD1	2.54	0.41
6:B:267:ARG:HG2	6:B:315:TYR:CZ	2.56	0.41
6:B:352:ASN:C	6:B:354:ARG:N	2.79	0.41
4:E:679:ARG:HB3	4:E:764:GLN:NE2	2.35	0.41
6:B:135:ILE:HD12	6:B:138:ARG:HH11	1.85	0.41
6:B:228:ARG:O	6:B:232:GLU:HG3	2.20	0.41
4:E:68:ASP:HB3	4:E:71:GLN:CD	2.45	0.41
4:E:842:TYR:CD2	4:E:843:SER:N	2.89	0.41
4:E:1158:ILE:O	4:E:1161:PRO:HD2	2.21	0.41
5:D:11:TRP:CD1	5:D:117:ASP:HA	2.56	0.41
4:E:259:ILE:HA	4:E:265:TYR:CE1	2.56	0.41
4:E:520:LEU:HD12	4:E:833:TYR:CE2	2.56	0.41
6:B:24:LEU:HA	6:B:24:LEU:HD23	1.88	0.41
6:B:136:MET:HE3	6:B:136:MET:HB2	1.84	0.41
6:B:172:VAL:HG13	6:B:229:GLU:HG2	2.02	0.41
4:E:780:ARG:O	4:E:784:PHE:CD2	2.73	0.41
6:B:311:LYS:HE3	6:B:311:LYS:HB2	1.91	0.41
5:D:58:ARG:HA	5:D:69:ARG:HE	1.85	0.40
6:B:203:LEU:HB3	6:B:252:LEU:HD12	2.03	0.40
4:E:346:PHE:HZ	6:B:111:ARG:NH1	2.19	0.40
4:E:1066:SER:O	4:E:1067:ILE:C	2.64	0.40
4:E:261:VAL:O	4:E:497:CYS:C	2.65	0.40
4:E:567:ARG:HH11	4:E:567:ARG:CG	2.29	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:1154:ILE:HG22	4:E:1159:THR:HG23	2.03	0.40
4:E:520:LEU:HB3	4:E:841:TRP:CZ3	2.57	0.40
4:E:653:ARG:NH2	4:E:768:PRO:HB3	2.37	0.40
4:E:657:ASP:OD2	4:E:657:ASP:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	24/26 (92%)	22 (92%)	2 (8%)	0	100	100
4	E	1116/2222 (50%)	1051 (94%)	52 (5%)	13 (1%)	10	41
5	D	130/132 (98%)	127 (98%)	3 (2%)	0	100	100
6	B	370/380 (97%)	342 (92%)	25 (7%)	3 (1%)	16	49
All	All	1640/2760 (59%)	1542 (94%)	82 (5%)	16 (1%)	15	45

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	E	833	TYR
4	E	845	GLU
4	E	846	MET
4	E	987	LYS
4	E	1053	SER
4	E	1055	THR
4	E	659	ILE
4	E	1061	GLY
6	B	342	GLU
4	E	128	SER
4	E	986	LEU

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Mol	Chain	Res	Type
6	B	340	ASP
6	B	353	SER
4	E	834	VAL
4	E	798	ILE
4	E	1043	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	20/24 (83%)	20 (100%)	0	100	100
4	E	950/2014 (47%)	944 (99%)	6 (1%)	78	79
5	D	110/119 (92%)	108 (98%)	2 (2%)	51	68
6	B	326/352 (93%)	325 (100%)	1 (0%)	86	83
All	All	1406/2509 (56%)	1397 (99%)	9 (1%)	76	79

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	E	167	LEU
4	E	795	LEU
4	E	815	LEU
4	E	825	VAL
4	E	827	LEU
4	E	834	VAL
5	D	67	LEU
5	D	73	LEU
6	B	28	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	728	ASN
4	E	39	ASN

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Mol	Chain	Res	Type
4	E	434	GLN
4	E	523	GLN
4	E	526	GLN
4	E	572	ASN
4	E	631	ASN
4	E	828	ASN
4	E	924	GLN
4	E	993	GLN
4	E	997	ASN
6	B	29	GLN
6	B	33	GLN
6	B	34	ASN
6	B	54	HIS
6	B	224	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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	SF4	E	2301	4	0,12,12	-	-	-		

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
7	SF4	E	2301	4	-	-	0/6/5/5

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.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	E	2301	SF4	6	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

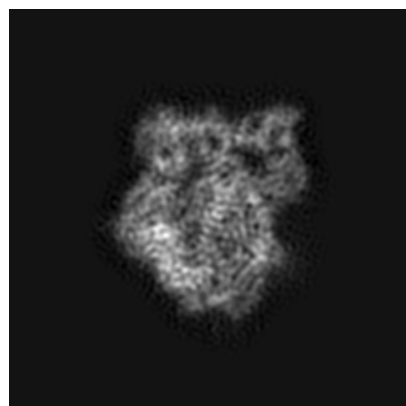
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41663. These allow visual inspection of the internal detail of the map and identification of artifacts.

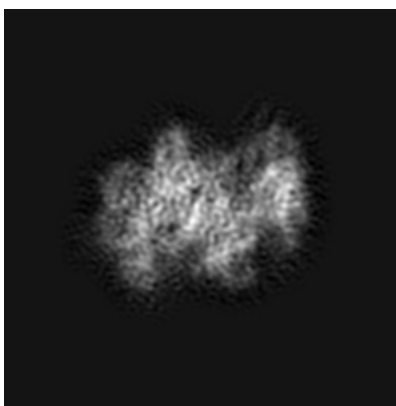
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

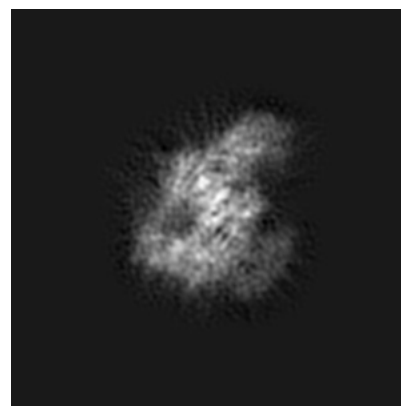
6.1.1 Primary map



X

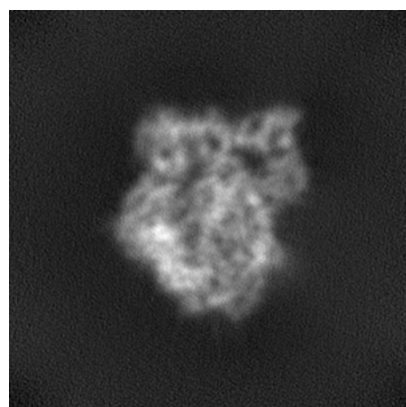


Y

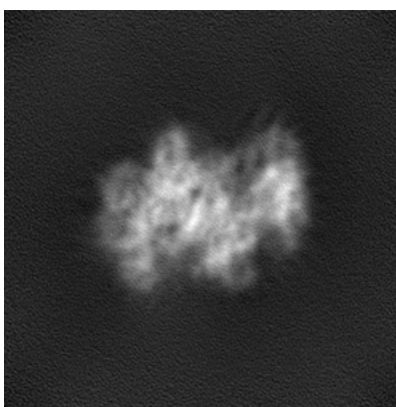


Z

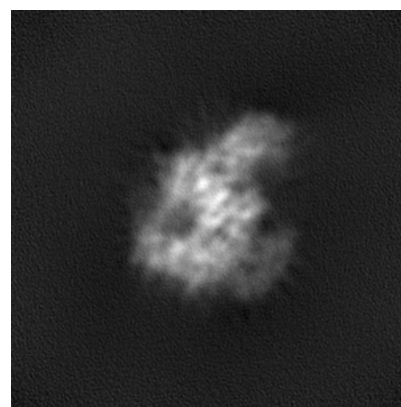
6.1.2 Raw map



X



Y

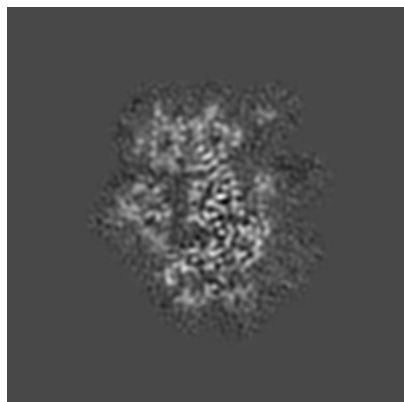


Z

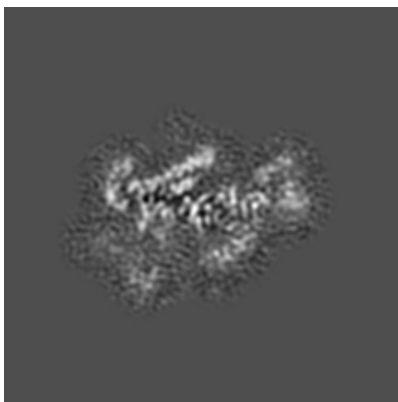
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

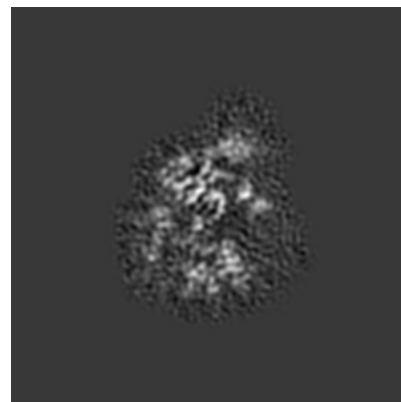
6.2.1 Primary map



X Index: 128

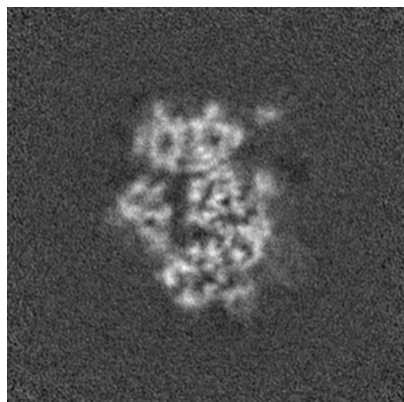


Y Index: 128

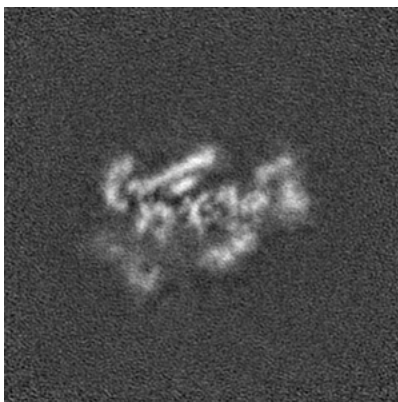


Z Index: 128

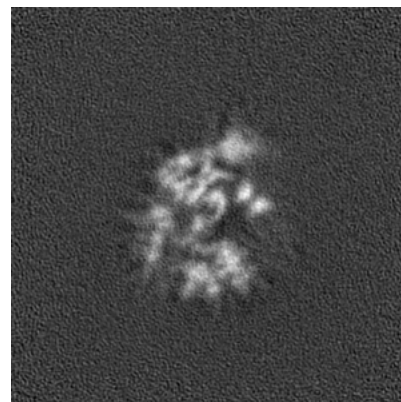
6.2.2 Raw map



X Index: 128



Y Index: 128

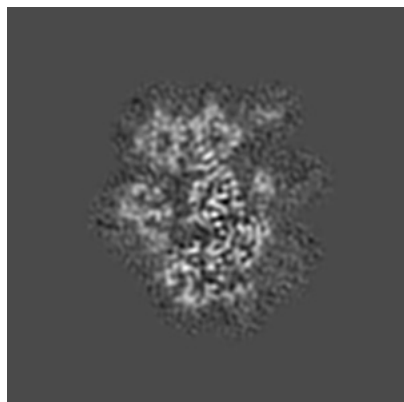


Z Index: 128

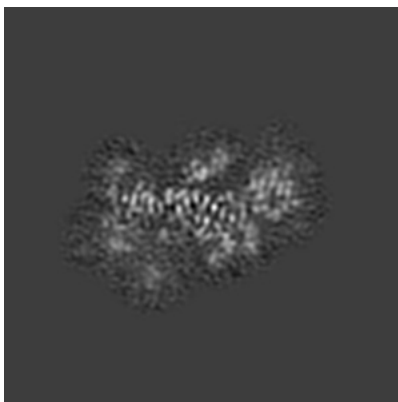
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

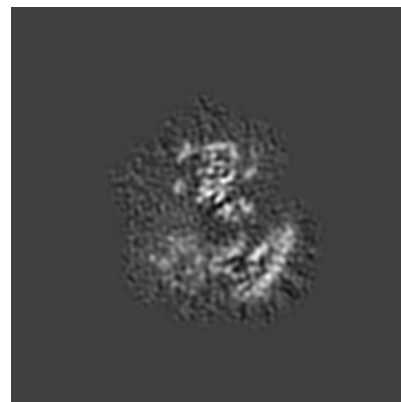
6.3.1 Primary map



X Index: 129

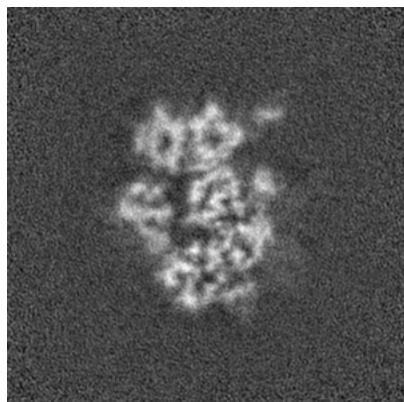


Y Index: 136

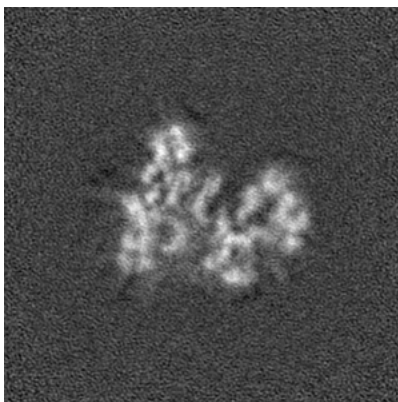


Z Index: 111

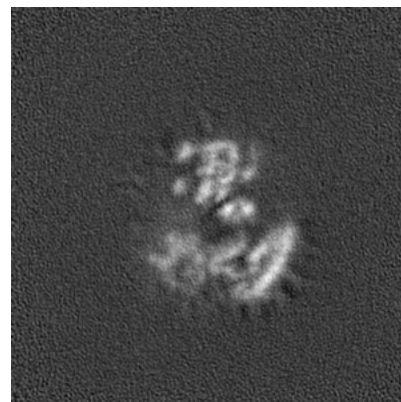
6.3.2 Raw map



X Index: 129



Y Index: 102

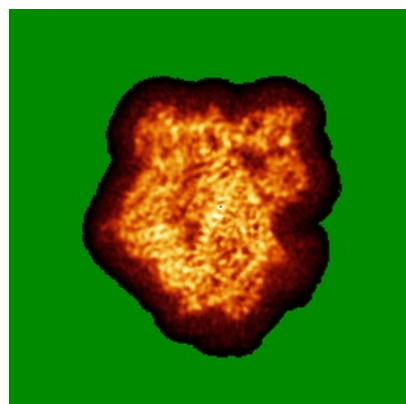


Z Index: 111

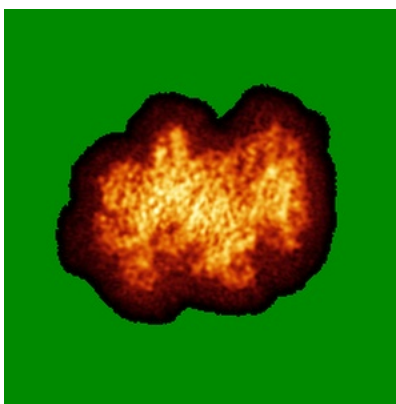
The images above show the largest variance slices of the map in three orthogonal directions.

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

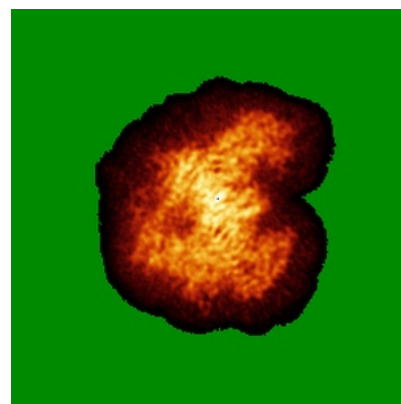
6.4.1 Primary map



X

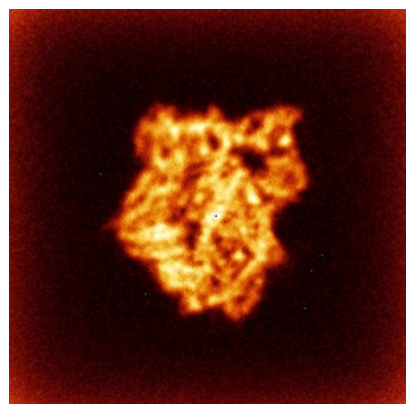


Y

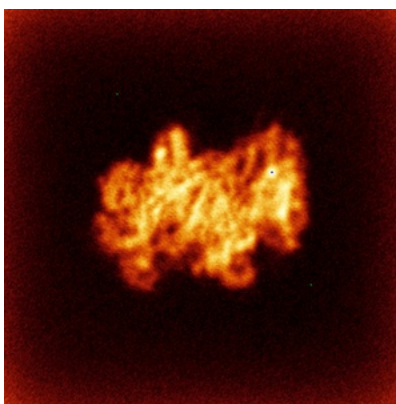


Z

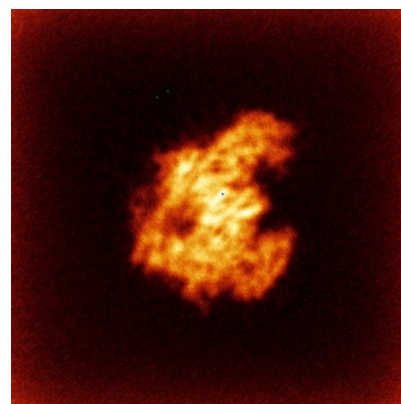
6.4.2 Raw map



X



Y

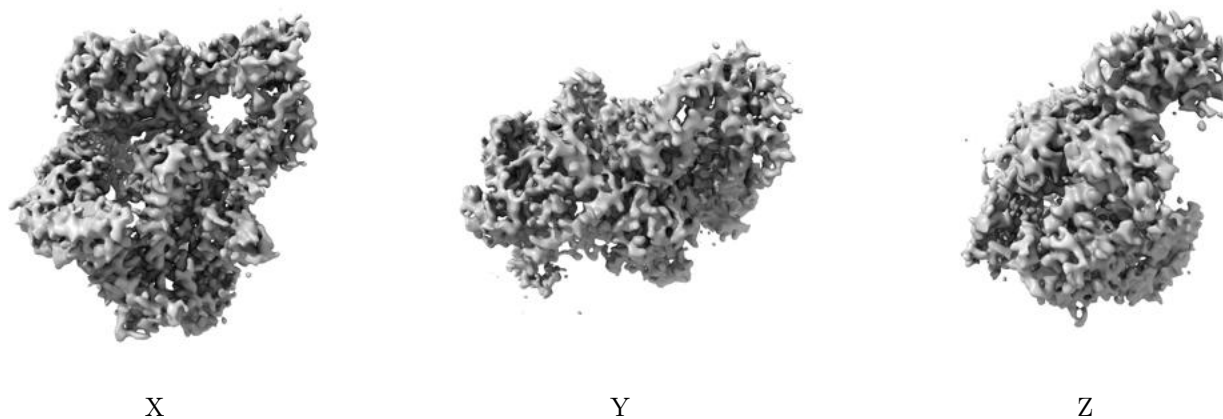


Z

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.

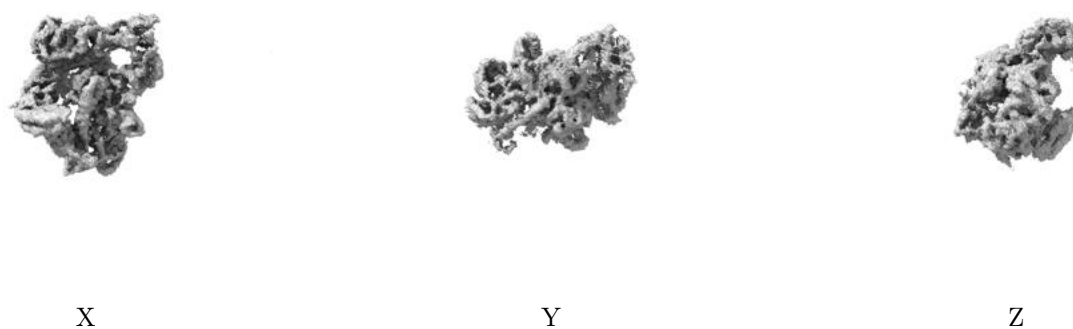
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.326. 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.

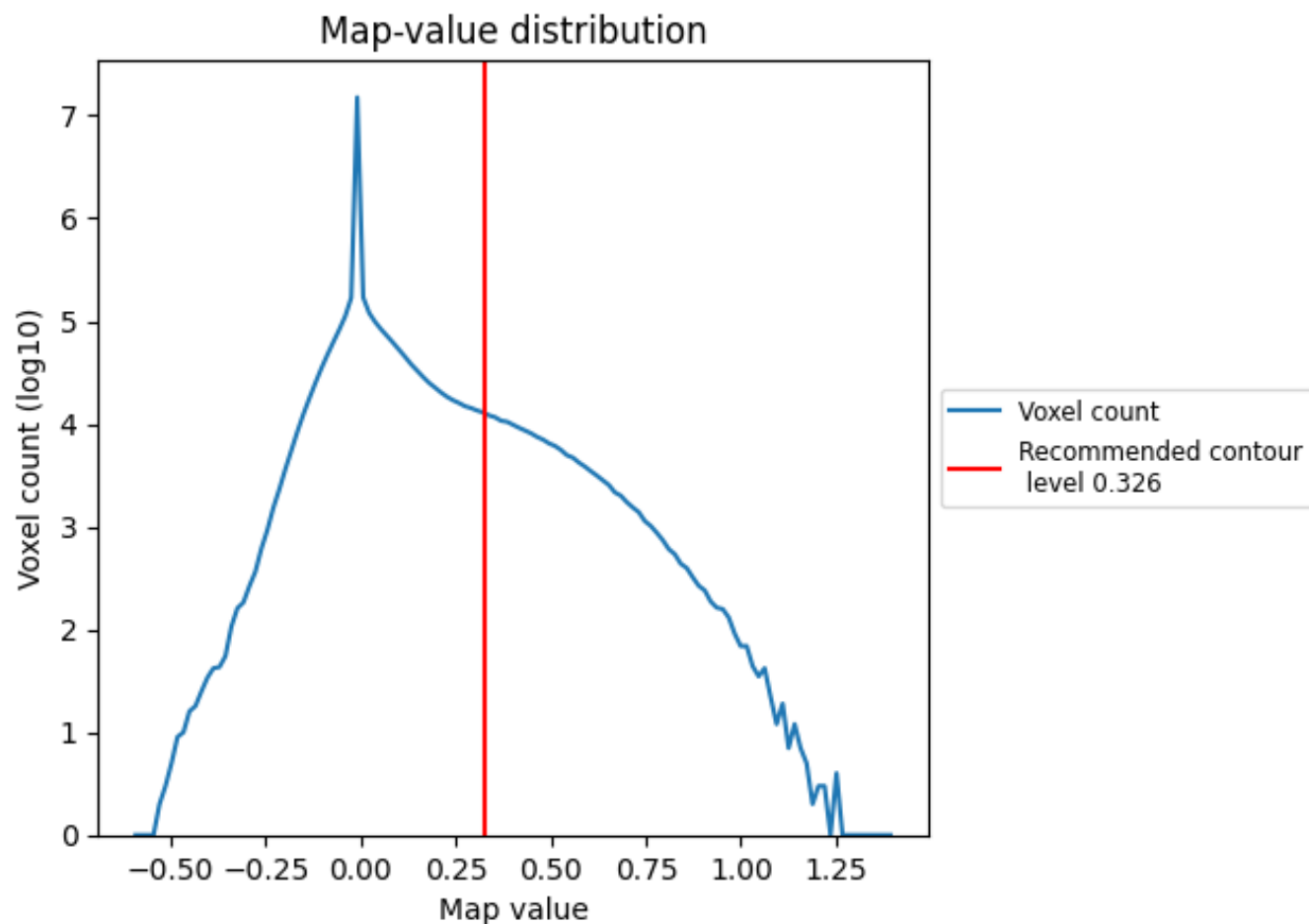
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

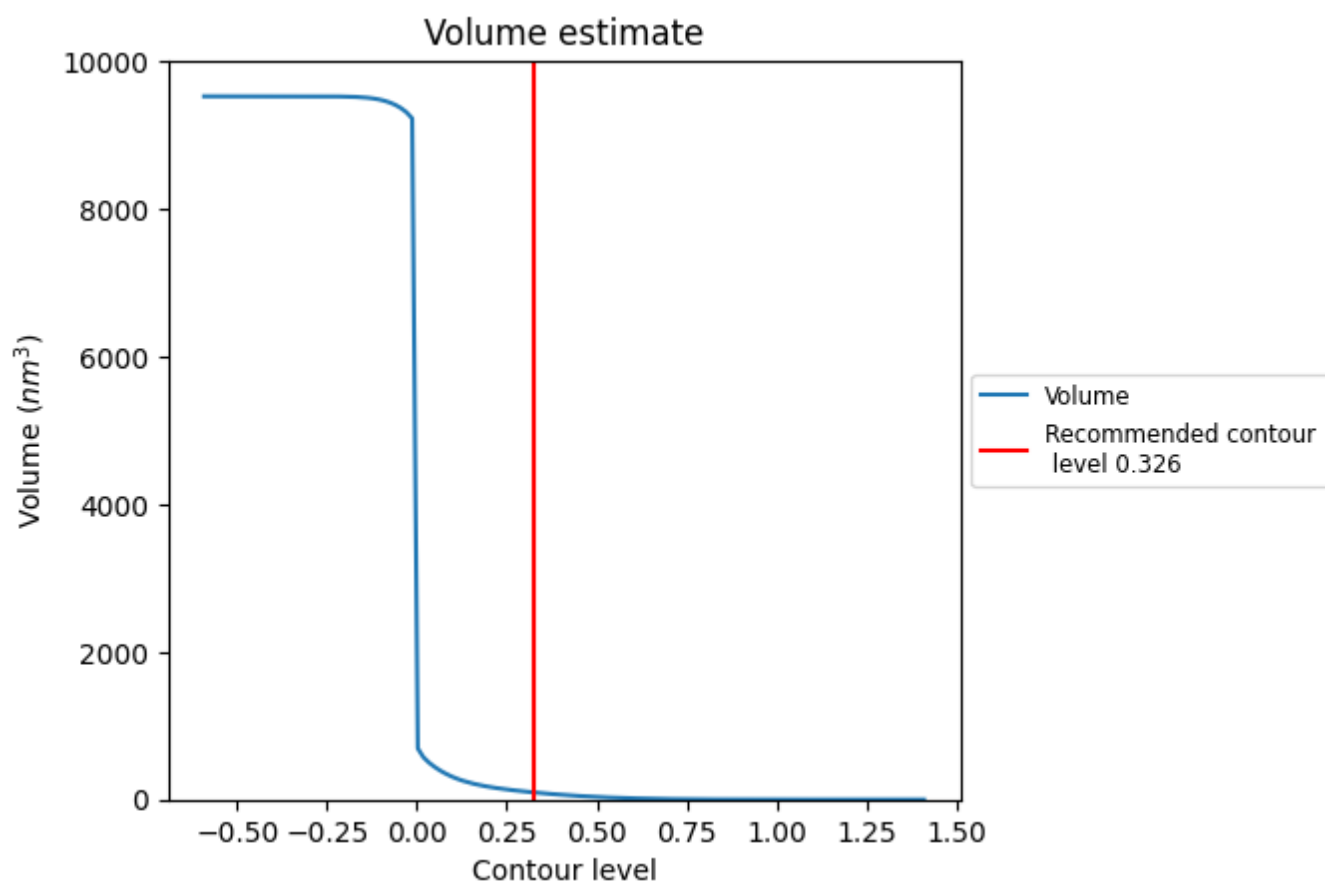
This section contains the results of statistical analysis of the map.

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.

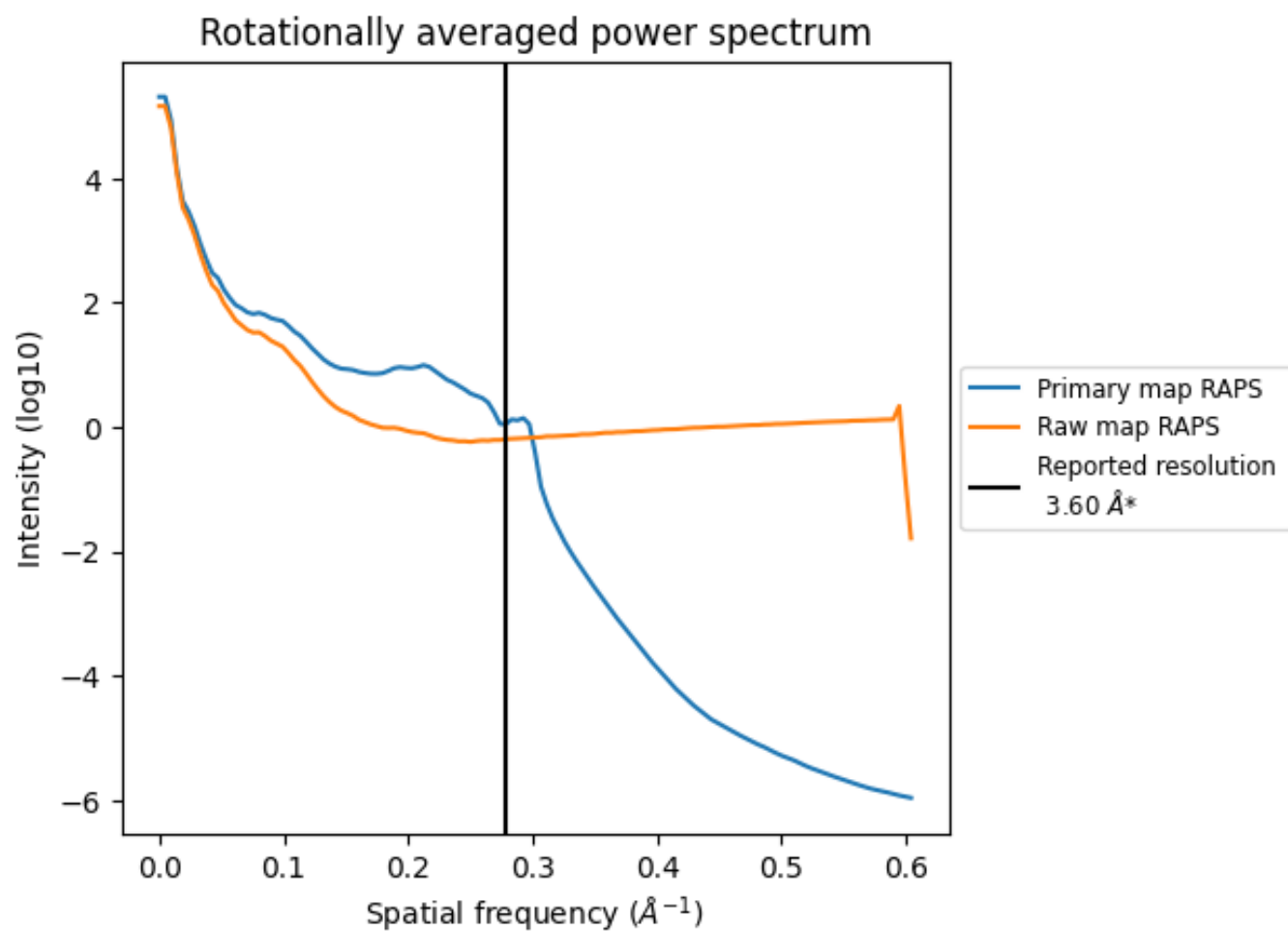
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 96 nm³; this corresponds to an approximate mass of 87 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 ⓘ

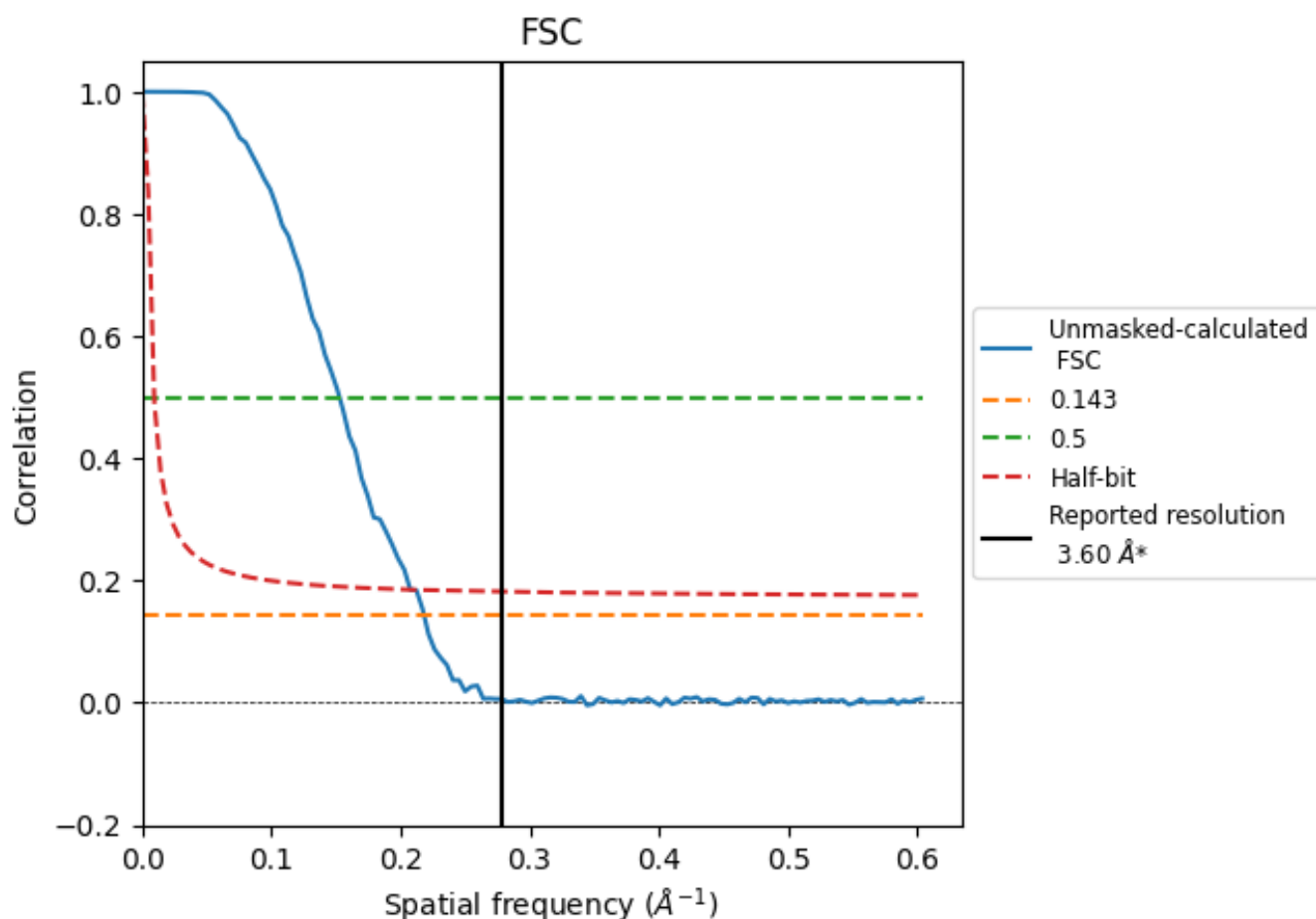


*Reported resolution corresponds to spatial frequency of 0.278 $\text{\AA}^{-1}$

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](#)

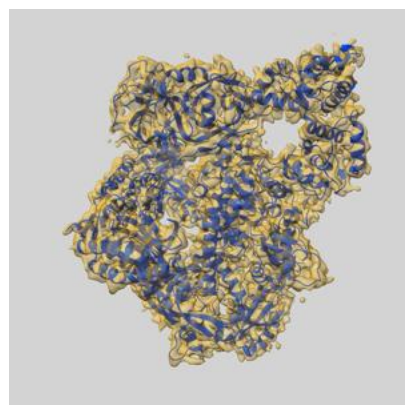
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.60	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.58	6.55	4.82

*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 4.58 differs from the reported value 3.6 by more than 10 %

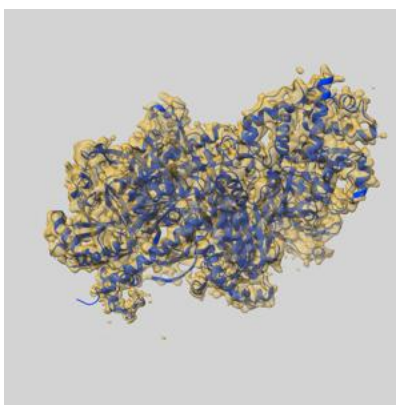
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-41663 and PDB model 8TW9. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

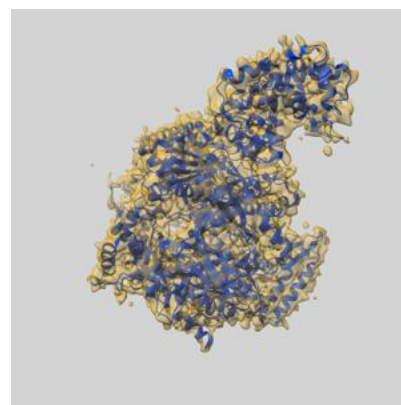
9.1 Map-model overlay [i](#)



X



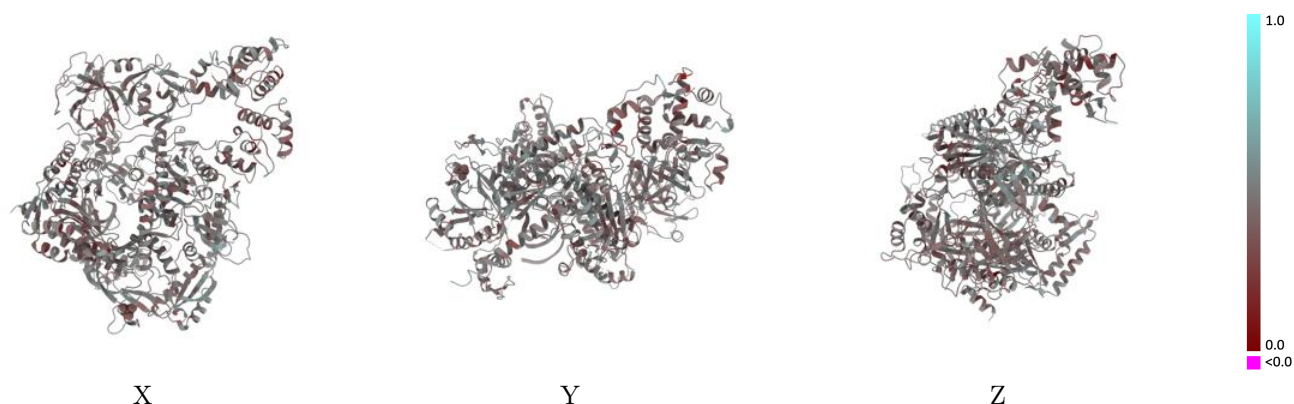
Y



Z

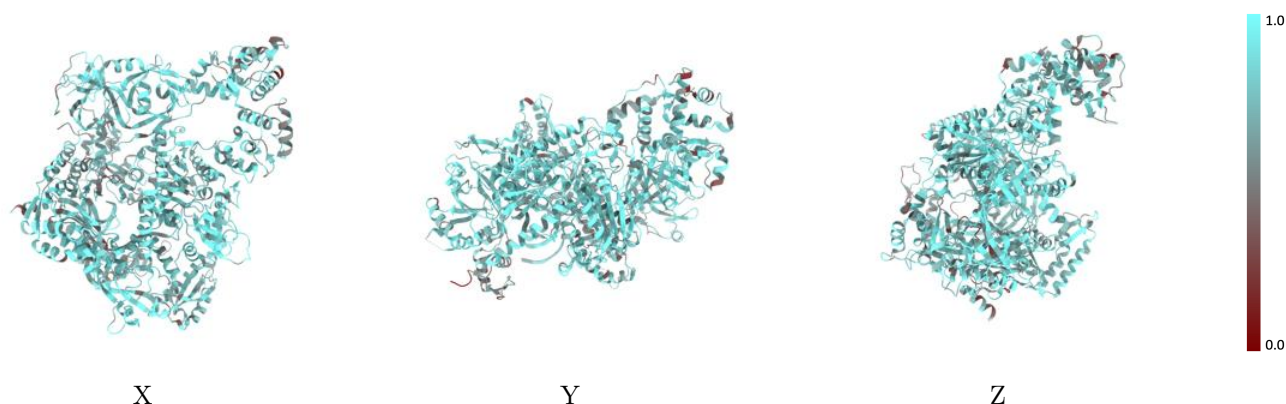
The images above show the 3D surface view of the map at the recommended contour level 0.326 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



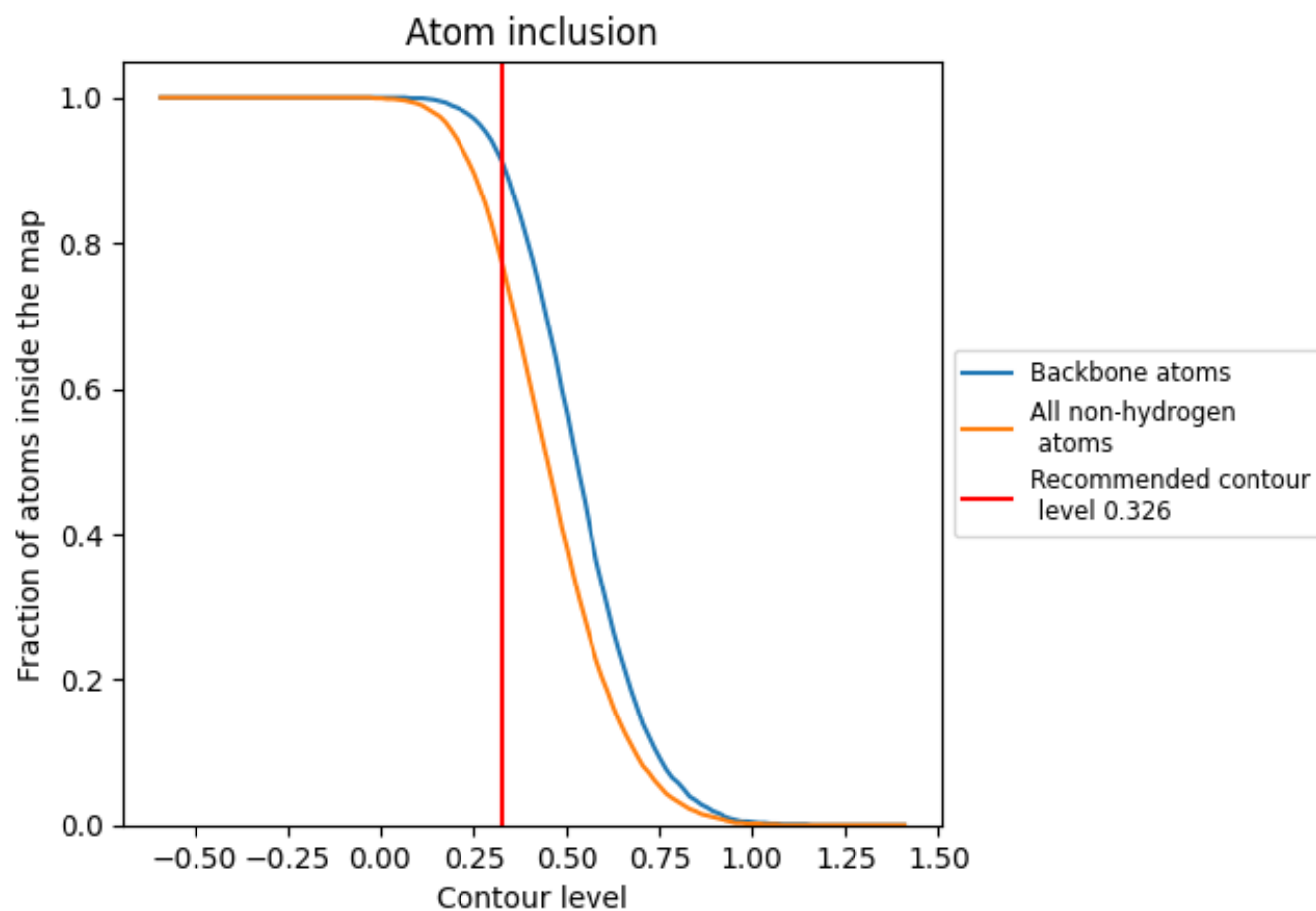
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.326).

9.4 Atom inclusion ⓘ



At the recommended contour level, 91% of all backbone atoms, 78% 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 (0.326) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7760	0.4380
B	0.7590	0.4310
C	0.7410	0.4140
D	0.7900	0.4270
E	0.7860	0.4430
P	0.7490	0.4260
T	0.6600	0.4240

1.0

0.0

<0.0