



Full wwPDB EM Validation Report ⓘ

Dec 1, 2025 – 04:28 pm GMT

PDB ID : 9RPW / pdb_00009rpw
EMDB ID : EMD-54173
Title : Cryo-EM structure of Shigella flexneri LptDE dimer: closed-state unbound and open-state bound by Oekolampad phage RBP
Authors : Dunbar, E.; Basle, A.; van den Berg, B.
Deposited on : 2025-06-25
Resolution : 3.20 Å(reported)
Based on initial model : .

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:

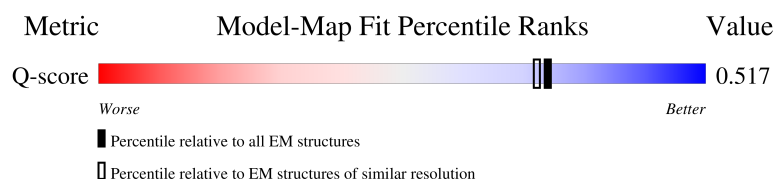
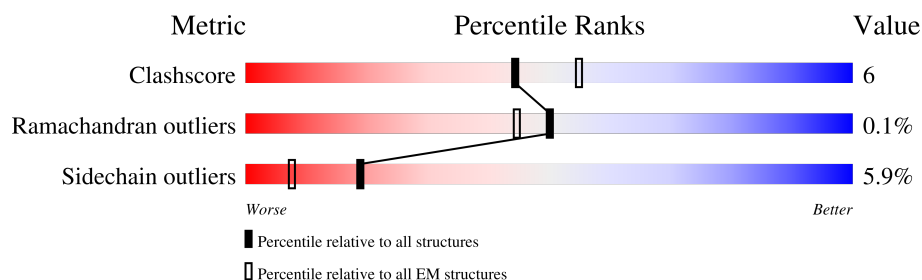
EMDB validation analysis : 0.0.1.dev129
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.46

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 3.20 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	210492	15764	-
Ramachandran outliers	207382	16835	-
Sidechain outliers	206894	16415	-
Q-score	-	25397	15020 (2.70 - 3.70)

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	784	
1	D	784	
2	B	199	
2	E	199	

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Mol	Chain	Length	Quality of chain
3	C	327	<div><div></div><div>65%</div><div>12%</div><div>•</div><div>22%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13588 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 LPS-assembly protein LptD.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	D	560	Total	C	N	O	S	0	0
			4594	2893	781	906	14		
1	A	572	Total	C	N	O	S	0	0
			4694	2960	799	921	14		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	initiating methionine	UNP Q83SQ0
D	2	GLU	-	expression tag	UNP Q83SQ0
A	1	MET	-	initiating methionine	UNP Q83SQ0
A	2	GLU	-	expression tag	UNP Q83SQ0

- Molecule 2 is a protein called LPS-assembly lipoprotein LptE.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	150	Total	C	N	O	S	0	0
			1181	737	215	221	8		
2	B	150	Total	C	N	O	S	0	0
			1181	737	215	221	8		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	194	GLU	-	expression tag	UNP Q83LX4
E	195	ASN	-	expression tag	UNP Q83LX4
E	196	LEU	-	expression tag	UNP Q83LX4
E	197	TYR	-	expression tag	UNP Q83LX4
E	198	PHE	-	expression tag	UNP Q83LX4
E	199	GLN	-	expression tag	UNP Q83LX4
B	194	GLU	-	expression tag	UNP Q83LX4
B	195	ASN	-	expression tag	UNP Q83LX4
B	196	LEU	-	expression tag	UNP Q83LX4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	197	TYR	-	expression tag	UNP Q83LX4
B	198	PHE	-	expression tag	UNP Q83LX4
B	199	GLN	-	expression tag	UNP Q83LX4

- Molecule 3 is a protein called Tail fiber protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	254	Total	C	N	O	S	0	0
			1938	1217	347	364	10		

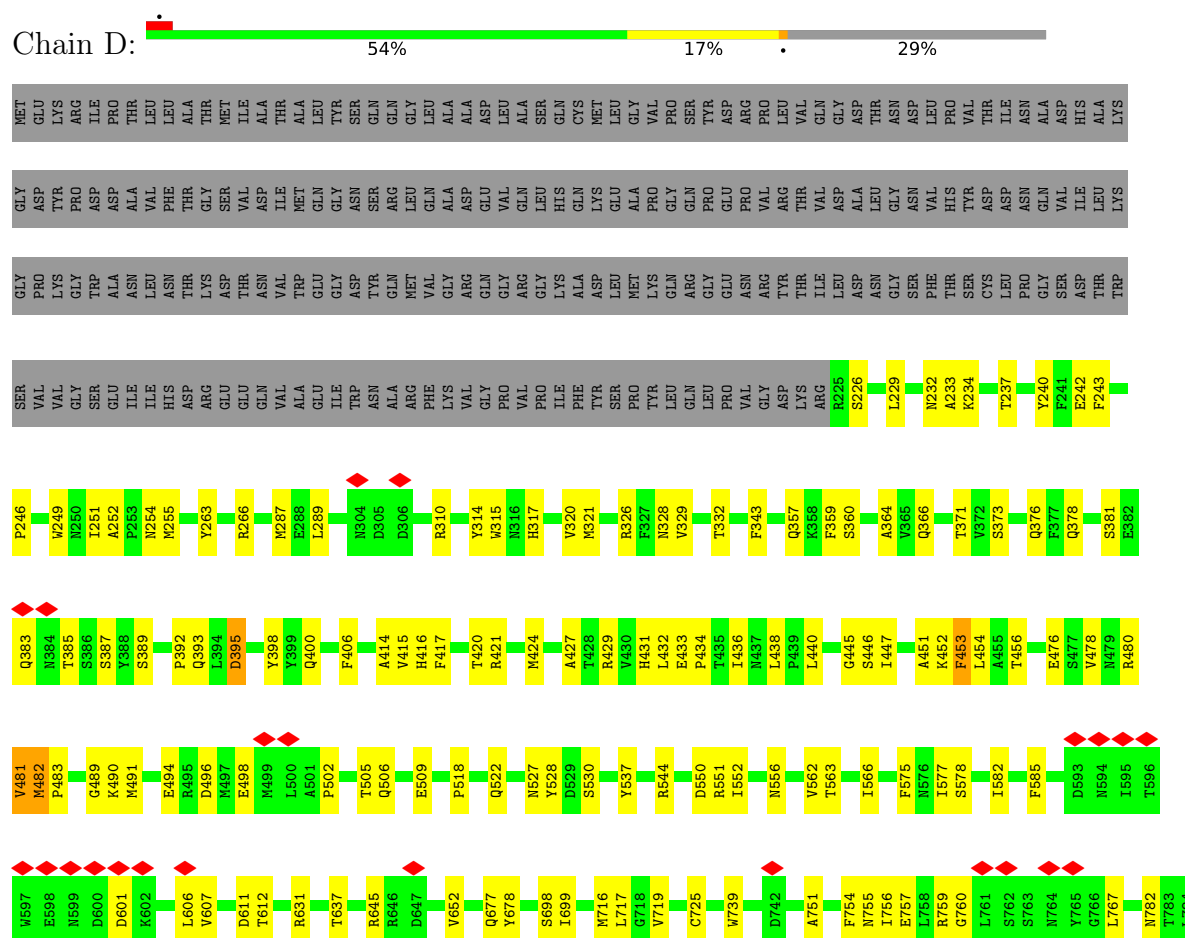
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	322	HIS	-	expression tag	UNP A0AAE8B389
C	323	HIS	-	expression tag	UNP A0AAE8B389
C	324	HIS	-	expression tag	UNP A0AAE8B389
C	325	HIS	-	expression tag	UNP A0AAE8B389
C	326	HIS	-	expression tag	UNP A0AAE8B389
C	327	HIS	-	expression tag	UNP A0AAE8B389

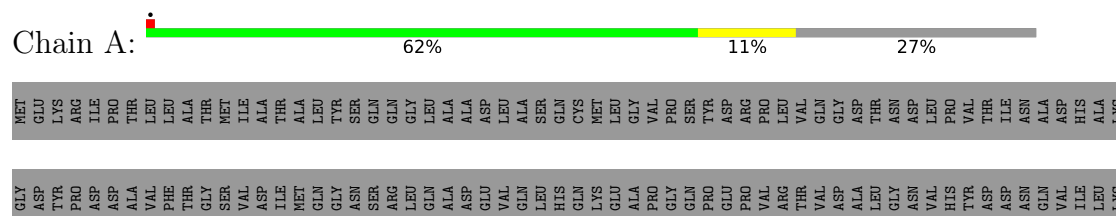
3 Residue-property plots

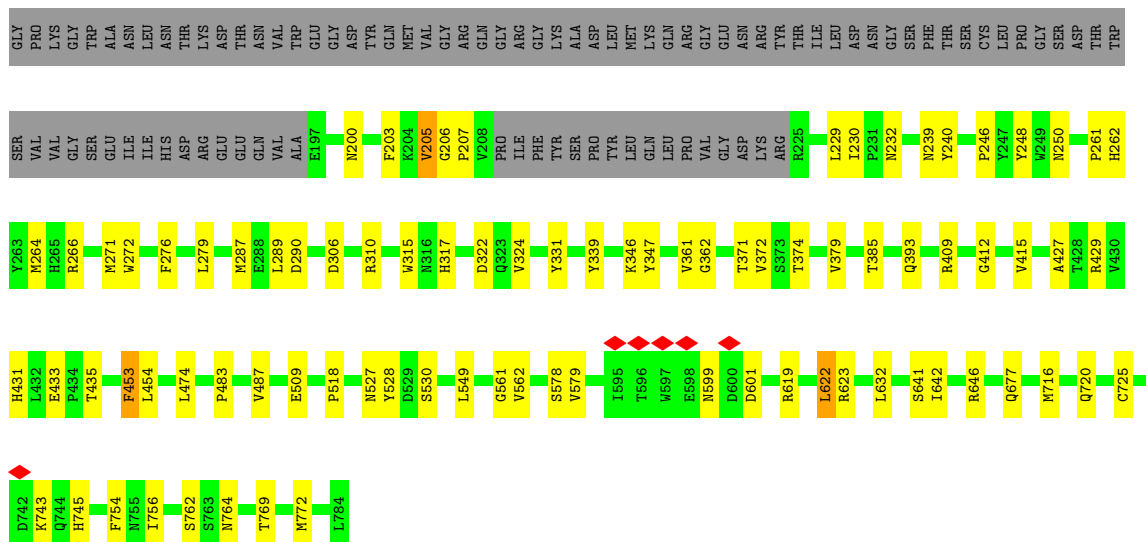
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: LPS-assembly protein LptD

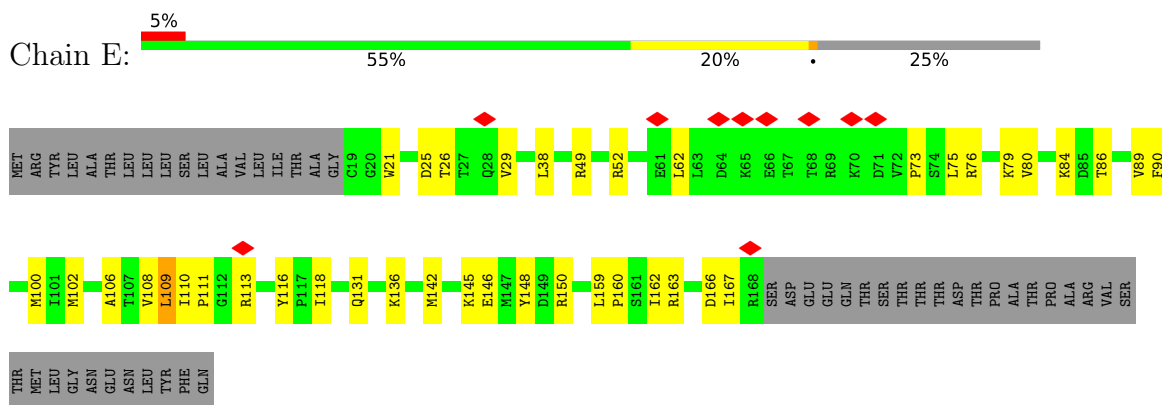


• Molecule 1: LPS-assembly protein LptD

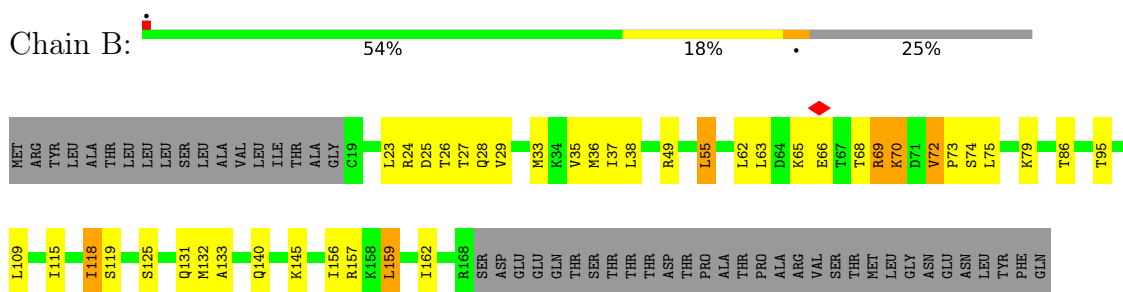




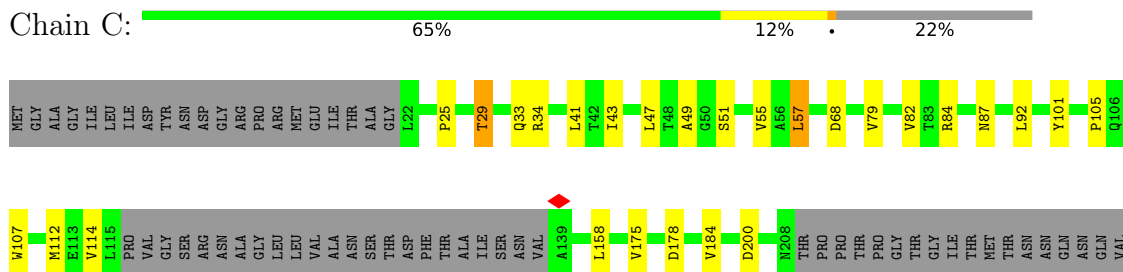
- Molecule 2: LPS-assembly lipoprotein LptE



- Molecule 2: LPS-assembly lipoprotein LptE



- Molecule 3: Tail fiber protein



THR	PHE	SER	THR	VAL	N232	K233	Y237	T245	L250	Q256	L257	L263	Y270	R274	M280	R291	S297	I317	M320	Y321	HIS	HIS	HIS	HIS	HIS	HIS
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	55494	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	35.6	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	165000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.484	Depositor
Minimum map value	-0.247	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.09	Depositor
Map size (\AA)	426.24, 426.24, 426.24	wwPDB
Map dimensions	576, 576, 576	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.74, 0.74, 0.74	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	0.27	0/4822	0.36	0/6561
1	D	0.22	0/4719	0.38	0/6422
2	B	0.27	0/1198	0.46	0/1617
2	E	0.24	0/1198	0.39	0/1617
3	C	0.27	0/1972	0.34	0/2681
All	All	0.25	0/13909	0.38	0/18898

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4694	0	4356	39	0
1	D	4594	0	4256	71	0
2	B	1181	0	1212	19	0
2	E	1181	0	1212	19	0
3	C	1938	0	1934	18	0
All	All	13588	0	12970	165	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 (165) 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:175:VAL:HG22	3:C:184:VAL:HG22	1.73	0.69
1:A:248:TYR:HE2	1:A:250:ASN:HB2	1.60	0.67
1:A:262:HIS:HB3	1:A:271:MET:HG2	1.77	0.67
3:C:33:GLN:HG3	3:C:41:LEU:HD11	1.78	0.65
1:D:498:GLU:HA	1:D:502:PRO:HA	1.78	0.64
2:B:29:VAL:HG12	2:B:159:LEU:HD23	1.81	0.62
1:D:251:ILE:HG13	1:D:252:ALA:H	1.64	0.61
1:D:421:ARG:HH21	1:D:424:MET:HE2	1.65	0.60
1:D:483:PRO:HD2	1:D:518:PRO:HG3	1.82	0.60
2:B:33:MET:HE1	2:B:162:ILE:HG13	1.85	0.59
1:D:328:ASN:HB2	1:D:360:SER:HB3	1.83	0.59
3:C:237:TYR:HA	3:C:317:ILE:HG22	1.84	0.59
1:D:582:ILE:HD11	1:D:607:VAL:HB	1.84	0.58
1:A:271:MET:HE2	1:A:290:ASP:HB3	1.87	0.57
1:A:361:VAL:HG12	1:A:362:GLY:H	1.68	0.57
1:D:677:GLN:HE21	1:D:739:TRP:CG	2.23	0.56
3:C:41:LEU:HD23	3:C:92:LEU:HD12	1.87	0.56
1:D:240:TYR:CE2	1:D:266:ARG:HD3	2.42	0.55
1:D:373:SER:HB2	1:D:393:GLN:HB3	1.87	0.55
1:D:453:PHE:HE1	1:D:481:VAL:HG12	1.71	0.55
1:D:544:ARG:HG2	1:D:552:ILE:HD12	1.87	0.55
2:E:163:ARG:HE	2:E:167:ILE:HD11	1.71	0.55
1:D:229:LEU:HB2	1:D:246:PRO:HG2	1.89	0.54
1:A:248:TYR:CE2	1:A:250:ASN:HB2	2.41	0.54
3:C:256:GLN:HG2	3:C:257:LEU:H	1.73	0.54
3:C:55:VAL:HG11	3:C:82:VAL:HG11	1.90	0.54
1:A:433:GLU:HG2	1:A:454:LEU:HD23	1.89	0.54
1:D:482:MET:HB2	1:D:518:PRO:HG2	1.89	0.54
1:D:601:ASP:HB2	1:D:631:ARG:HG3	1.90	0.53
2:B:36:MET:HA	2:B:72:VAL:HG13	1.91	0.53
1:D:530:SER:HB2	1:D:550:ASP:HB2	1.91	0.53
1:D:506:GLN:HB2	1:D:566:ILE:HG22	1.92	0.52
1:A:206:GLY:HA3	1:A:232:ASN:HB3	1.91	0.52
1:A:315:TRP:HE1	1:A:317:HIS:HB2	1.72	0.52
1:D:480:ARG:HH11	1:D:482:MET:HE1	1.74	0.52
2:B:69:ARG:HG2	2:B:70:LYS:H	1.75	0.52
1:A:427:ALA:HB2	1:A:528:TYR:CZ	2.45	0.52
1:D:645:ARG:HH21	1:D:782:ASN:HB3	1.76	0.51
1:D:527:ASN:HB2	1:D:551:ARG:HH21	1.74	0.51
1:D:432:LEU:O	1:D:434:PRO:HD3	2.11	0.51
1:A:315:TRP:NE1	1:A:317:HIS:HB2	2.27	0.50
3:C:25:PRO:HB2	3:C:112:MET:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:577:ILE:HG13	1:D:612:THR:HB	1.94	0.50
1:D:417:PHE:HD2	1:D:427:ALA:HB3	1.76	0.50
2:E:38:LEU:HD23	2:E:52:ARG:HE	1.78	0.49
1:D:677:GLN:HE21	1:D:739:TRP:CD1	2.31	0.48
1:D:754:PHE:CE2	1:D:756:ILE:HD11	2.49	0.48
2:E:21:TRP:CD1	2:E:21:TRP:H	2.31	0.48
2:B:36:MET:HB3	2:B:73:PRO:HG2	1.95	0.48
3:C:49:ALA:HA	3:C:87:ASN:HD21	1.78	0.48
3:C:233:LYS:HD3	3:C:320:MET:HE2	1.95	0.48
2:E:113:ARG:HH12	2:E:167:ILE:HD13	1.79	0.48
1:A:346:LYS:HD2	1:A:347:TYR:CE2	2.49	0.48
1:A:409:ARG:HG2	1:A:435:THR:HG22	1.95	0.48
1:A:720:GLN:HB2	1:A:772:MET:HE1	1.96	0.47
2:E:113:ARG:HD3	2:E:166:ASP:CG	2.39	0.47
2:B:69:ARG:HG2	2:B:70:LYS:N	2.29	0.47
2:B:28:GLN:HG2	2:B:29:VAL:H	1.79	0.47
1:D:436:ILE:HG22	1:D:451:ALA:O	2.14	0.47
1:D:315:TRP:HE1	1:D:317:HIS:HB2	1.80	0.47
3:C:29:THR:HG21	3:C:43:ILE:HD11	1.97	0.47
1:D:254:ASN:OD1	1:D:255:MET:HG3	2.15	0.47
2:E:76:ARG:HH12	2:E:109:LEU:HD11	1.80	0.46
2:B:118:ILE:HG22	2:B:119:SER:H	1.80	0.46
2:B:55:LEU:HD13	2:B:156:ILE:HD12	1.96	0.46
1:D:376:GLN:HG2	1:D:389:SER:HB3	1.96	0.46
1:D:242:GLU:HA	1:D:263:TYR:O	2.16	0.46
1:D:431:HIS:CD2	1:D:433:GLU:HG3	2.50	0.46
3:C:250:ILE:HD11	3:C:280:MET:HE2	1.97	0.46
2:B:157:ARG:C	2:B:159:LEU:H	2.24	0.46
3:C:51:SER:O	3:C:114:VAL:HG12	2.16	0.46
3:C:105:PRO:HG2	3:C:107:TRP:CZ2	2.51	0.45
1:D:415:VAL:HG22	1:D:416:HIS:H	1.80	0.45
1:A:415:VAL:HG12	1:A:429:ARG:HB3	1.99	0.45
1:D:606:LEU:HD23	1:D:631:ARG:HH21	1.81	0.45
1:D:232:ASN:HD22	1:D:757:GLU:HA	1.82	0.45
1:D:522:GLN:HE21	1:D:522:GLN:HB2	1.64	0.45
2:E:146:GLU:O	2:E:150:ARG:HD2	2.16	0.45
1:A:207:PRO:HG3	1:A:230:ILE:HB	1.99	0.45
1:D:406:PHE:CE1	1:D:438:LEU:HD13	2.52	0.45
1:D:427:ALA:HB2	1:D:528:TYR:CZ	2.51	0.45
1:D:251:ILE:CG2	1:D:255:MET:HB2	2.47	0.45
1:D:234:LYS:HB3	1:D:755:ASN:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:PHE:HD1	1:A:754:PHE:HB3	1.82	0.44
3:C:57:LEU:HD23	3:C:79:VAL:HG11	1.98	0.44
1:A:453:PHE:HD1	1:A:453:PHE:HA	1.72	0.44
2:B:132:MET:HG2	2:B:133:ALA:H	1.83	0.44
1:A:261:PRO:HB3	1:A:272:TRP:CZ3	2.52	0.44
1:A:622:LEU:HD23	1:A:623:ARG:N	2.33	0.44
2:E:113:ARG:HD3	2:E:166:ASP:OD1	2.18	0.44
1:D:480:ARG:NH1	1:D:482:MET:HE1	2.33	0.44
1:D:491:MET:O	1:D:509:GLU:HA	2.17	0.44
1:A:276:PHE:HB3	1:A:287:MET:HB2	1.99	0.44
2:E:136:LYS:HD2	2:E:136:LYS:HA	1.84	0.44
1:D:357:GLN:NE2	1:D:378:GLN:HB2	2.33	0.43
1:D:440:LEU:HB2	1:D:447:ILE:O	2.18	0.43
1:D:329:VAL:HG22	1:D:359:PHE:HD1	1.82	0.43
2:B:145:LYS:HA	2:B:145:LYS:HD2	1.74	0.43
1:D:453:PHE:CE1	1:D:481:VAL:HG12	2.53	0.43
1:A:262:HIS:CE1	1:A:264:MET:HE2	2.54	0.43
1:A:716:MET:HE3	1:A:716:MET:HB3	1.96	0.43
1:D:326:ARG:HH12	1:D:364:ALA:HB3	1.82	0.43
2:B:24:ARG:HD2	2:B:27:THR:HB	2.01	0.43
1:D:310:ARG:HE	1:D:343:PHE:HE2	1.66	0.43
1:D:556:ASN:O	1:D:585:PHE:HB2	2.19	0.43
1:A:622:LEU:HD23	1:A:623:ARG:H	1.83	0.43
1:D:233:ALA:HB2	1:D:243:PHE:HD1	1.84	0.43
2:E:142:MET:HE3	2:E:142:MET:HB2	1.91	0.43
3:C:263:LEU:HD22	3:C:274:ARG:HB2	2.01	0.43
2:E:89:VAL:HG22	2:E:90:PHE:O	2.19	0.42
2:B:37:ILE:HG23	2:B:63:LEU:HD23	2.01	0.42
2:B:86:THR:HG23	2:B:95:THR:HG23	2.01	0.42
1:D:371:THR:HB	1:D:395:ASP:OD1	2.19	0.42
1:A:412:GLY:HA2	1:A:431:HIS:O	2.20	0.42
1:D:716:MET:HE3	1:D:716:MET:HB3	1.86	0.42
1:D:725:CYS:HB3	1:A:725:CYS:HB2	2.00	0.42
2:E:102:MET:HE1	2:E:148:TYR:HD1	1.84	0.42
2:E:73:PRO:HA	2:E:111:PRO:HD3	2.00	0.42
3:C:270:TYR:HD1	3:C:297:SER:HA	1.84	0.42
1:A:205:VAL:HB	1:A:756:ILE:HB	2.02	0.42
2:B:74:SER:OG	2:B:109:LEU:HB3	2.20	0.42
1:D:249:TRP:CZ3	1:D:251:ILE:HB	2.55	0.42
1:A:483:PRO:HD2	1:A:518:PRO:HG3	2.01	0.42
1:A:562:VAL:O	1:A:578:SER:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:619:ARG:HD2	1:A:646:ARG:HD3	2.00	0.42
3:C:68:ASP:HB2	3:C:101:TYR:HE1	1.84	0.42
2:E:116:TYR:HB3	2:E:162:ILE:HD12	2.00	0.42
1:D:759:ARG:HG3	1:D:760:GLY:H	1.85	0.42
2:B:26:THR:C	2:B:28:GLN:H	2.27	0.42
1:D:414:ALA:HA	1:D:429:ARG:O	2.20	0.41
2:E:84:LYS:HB3	2:E:100:MET:HG2	2.02	0.41
2:E:159:LEU:HB2	2:E:160:PRO:HD3	2.01	0.41
1:D:237:THR:HG21	1:D:751:ALA:HA	2.01	0.41
1:D:387:SER:H	1:D:420:THR:HG22	1.85	0.41
1:D:496:ASP:HB3	1:D:498:GLU:CD	2.44	0.41
3:C:47:LEU:HD23	3:C:84:ARG:NE	2.35	0.41
1:D:381:SER:OG	1:D:383:GLN:HG2	2.20	0.41
1:D:446:SER:O	1:D:489:GLY:HA2	2.21	0.41
1:A:599:ASN:ND2	1:A:601:ASP:HB2	2.36	0.41
1:D:699:ILE:HG22	1:D:719:VAL:HG22	2.03	0.41
1:A:239:ASN:HA	1:A:266:ARG:HD3	2.03	0.41
1:D:456:THR:HB	1:D:480:ARG:HB3	2.03	0.41
2:E:106:ALA:HB3	2:E:118:ILE:HG13	2.02	0.41
1:A:743:LYS:HE2	1:A:745:HIS:NE2	2.36	0.41
1:D:445:GLY:HA2	1:D:490:LYS:O	2.21	0.41
2:E:102:MET:HE1	2:E:148:TYR:CD1	2.56	0.41
2:B:55:LEU:HD12	2:B:55:LEU:HA	1.80	0.41
1:D:566:ILE:CG1	1:D:575:PHE:HB3	2.51	0.41
1:D:698:SER:OG	1:D:767:LEU:HB3	2.21	0.41
2:E:38:LEU:HD12	2:E:38:LEU:HA	1.93	0.41
1:D:398:TYR:HE1	1:D:400:GLN:HB2	1.85	0.40
1:A:310:ARG:HH21	1:A:339:TYR:HA	1.87	0.40
1:A:393:GLN:HG3	1:A:415:VAL:HG23	2.02	0.40
1:A:561:GLY:HA2	1:A:579:VAL:O	2.22	0.40
1:D:537:TYR:CZ	1:D:578:SER:HB3	2.56	0.40
1:D:476:GLU:O	1:D:478:VAL:HG23	2.21	0.40
1:A:527:ASN:OD1	1:A:530:SER:HB2	2.20	0.40
2:B:70:LYS:HA	2:B:70:LYS:HD3	1.83	0.40
1:D:434:PRO:O	1:D:452:LYS:HB2	2.20	0.40
1:A:322:ASP:O	1:A:324:VAL:HG22	2.21	0.40
1:D:287:MET:HA	1:D:314:TYR:O	2.21	0.40
1:D:385:THR:OG1	1:D:421:ARG:HD3	2.22	0.40
1:D:453:PHE:CE2	1:D:483:PRO:HB3	2.57	0.40
1:A:229:LEU:HB2	1:A:246:PRO:HG2	2.04	0.40
1:A:762:SER:HB2	1:A:764:ASN:OD1	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	568/784 (72%)	494 (87%)	74 (13%)	0	100	100
1	D	558/784 (71%)	484 (87%)	73 (13%)	1 (0%)	44	75
2	B	148/199 (74%)	130 (88%)	17 (12%)	1 (1%)	19	54
2	E	148/199 (74%)	137 (93%)	11 (7%)	0	100	100
3	C	248/327 (76%)	220 (89%)	28 (11%)	0	100	100
All	All	1670/2293 (73%)	1465 (88%)	203 (12%)	2 (0%)	50	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	392	PRO
2	B	23	LEU

5.3.2 Protein sidechains [i](#)

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	503/681 (74%)	480 (95%)	23 (5%)	23	56
1	D	493/681 (72%)	473 (96%)	20 (4%)	26	59
2	B	131/173 (76%)	111 (85%)	20 (15%)	2	11
2	E	131/173 (76%)	117 (89%)	14 (11%)	5	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	C	212/272 (78%)	203 (96%)	9 (4%)	25 58
All	All	1470/1980 (74%)	1384 (94%)	86 (6%)	19 48

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	226	SER
1	D	289	LEU
1	D	320	VAL
1	D	321	MET
1	D	332	THR
1	D	366	GLN
1	D	395	ASP
1	D	453	PHE
1	D	454	LEU
1	D	481	VAL
1	D	482	MET
1	D	494	GLU
1	D	505	THR
1	D	562	VAL
1	D	563	THR
1	D	611	ASP
1	D	637	THR
1	D	652	VAL
1	D	678	TYR
1	D	717	LEU
2	E	25	ASP
2	E	26	THR
2	E	29	VAL
2	E	49	ARG
2	E	62	LEU
2	E	75	LEU
2	E	79	LYS
2	E	80	VAL
2	E	86	THR
2	E	108	VAL
2	E	109	LEU
2	E	110	ILE
2	E	131	GLN
2	E	145	LYS
1	A	200	ASN
1	A	205	VAL

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Mol	Chain	Res	Type
1	A	240	TYR
1	A	279	LEU
1	A	289	LEU
1	A	306	ASP
1	A	331	TYR
1	A	371	THR
1	A	372	VAL
1	A	374	THR
1	A	379	VAL
1	A	385	THR
1	A	453	PHE
1	A	474	LEU
1	A	487	VAL
1	A	509	GLU
1	A	549	LEU
1	A	622	LEU
1	A	632	LEU
1	A	641	SER
1	A	642	ILE
1	A	677	GLN
1	A	769	THR
2	B	25	ASP
2	B	35	VAL
2	B	38	LEU
2	B	49	ARG
2	B	55	LEU
2	B	62	LEU
2	B	65	LYS
2	B	66	GLU
2	B	68	THR
2	B	69	ARG
2	B	70	LYS
2	B	72	VAL
2	B	75	LEU
2	B	79	LYS
2	B	115	ILE
2	B	118	ILE
2	B	125	SER
2	B	131	GLN
2	B	140	GLN
2	B	159	LEU
3	C	29	THR

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Mol	Chain	Res	Type
3	C	34	ARG
3	C	57	LEU
3	C	158	LEU
3	C	178	ASP
3	C	200	ASP
3	C	245	THR
3	C	263	LEU
3	C	291	ARG

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

Mol	Chain	Res	Type
1	D	232	ASN
1	D	323	GLN
1	D	594	ASN
1	D	745	HIS
2	E	99	GLN
1	A	367	ASN
1	A	384	ASN
1	A	393	GLN
1	A	401	ASN
1	A	470	ASN
1	A	557	GLN
1	A	639	ASN
1	A	680	ASN
2	B	22	HIS
2	B	99	GLN
3	C	207	ASN
3	C	242	ASN

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 [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.

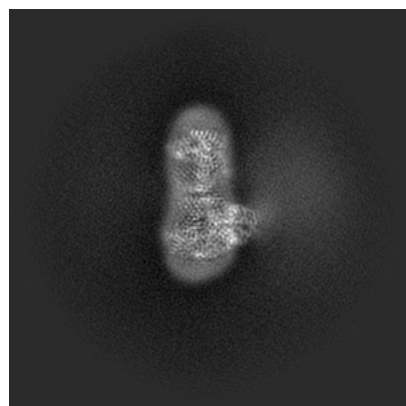
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-54173. These allow visual inspection of the internal detail of the map and identification of artifacts.

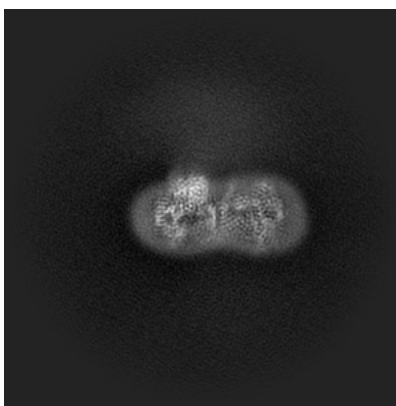
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

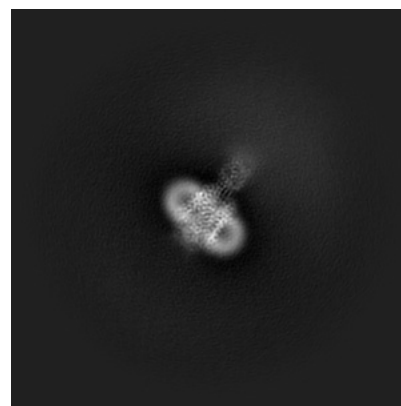
6.1.1 Primary map



X

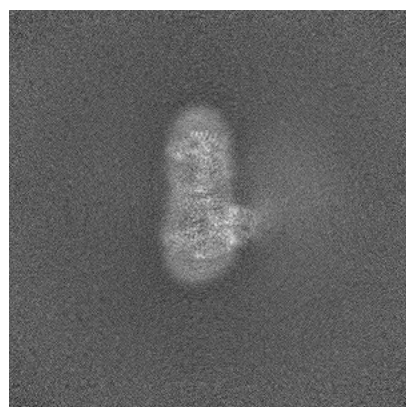


Y

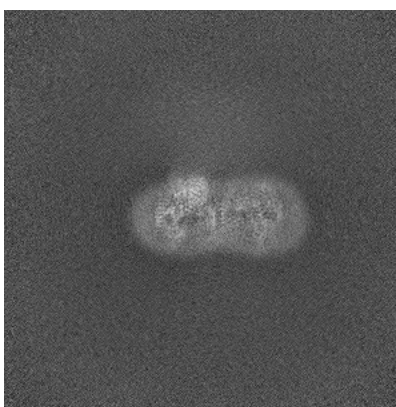


Z

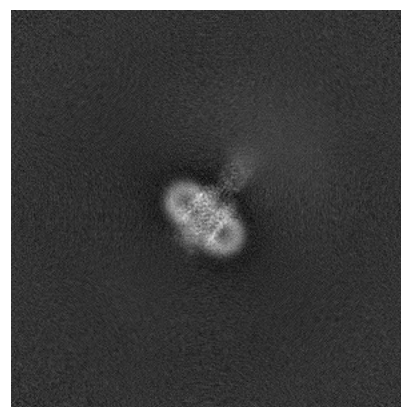
6.1.2 Raw map



X



Y

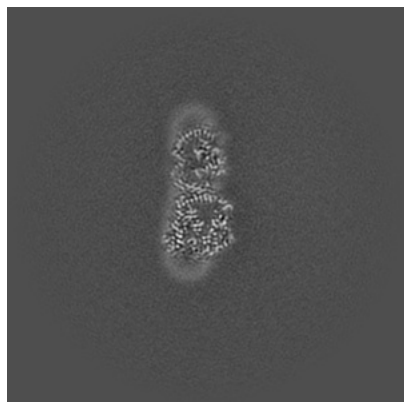


Z

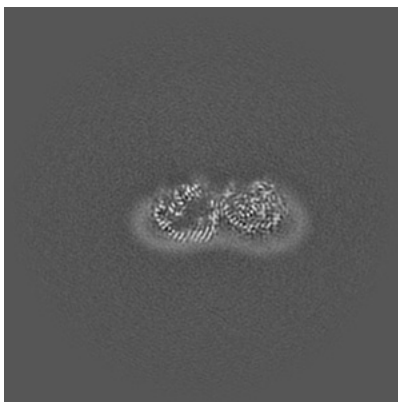
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

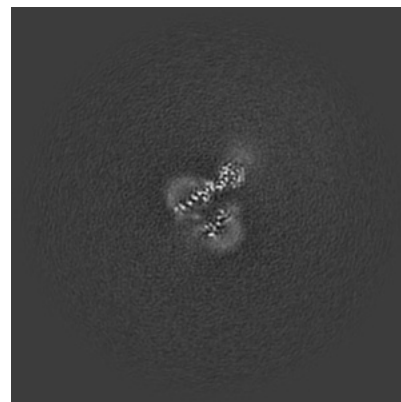
6.2.1 Primary map



X Index: 288

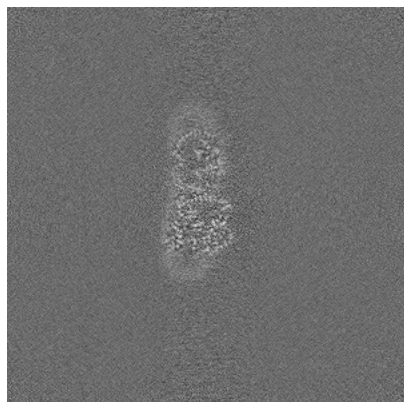


Y Index: 288

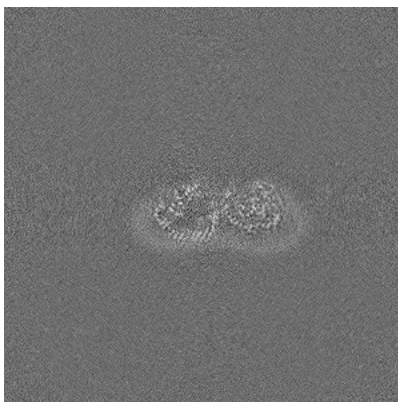


Z Index: 288

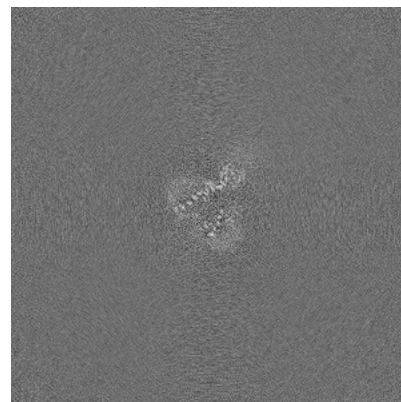
6.2.2 Raw map



X Index: 288



Y Index: 288

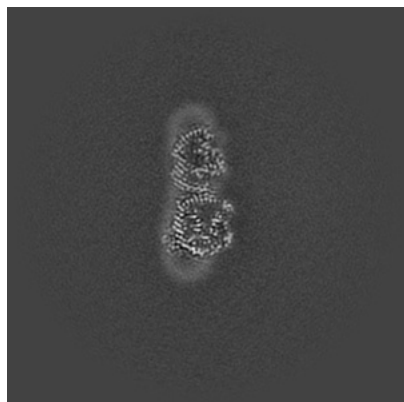


Z Index: 288

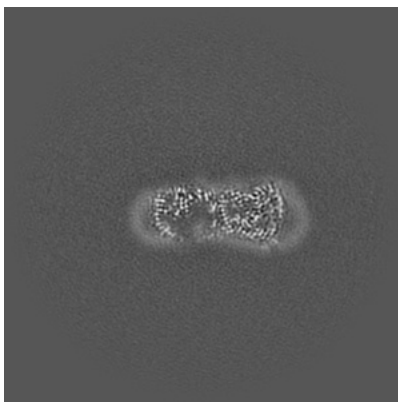
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

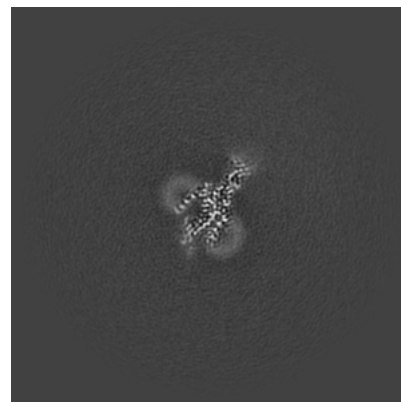
6.3.1 Primary map



X Index: 290

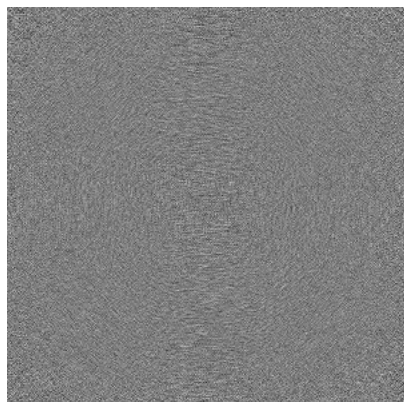


Y Index: 278

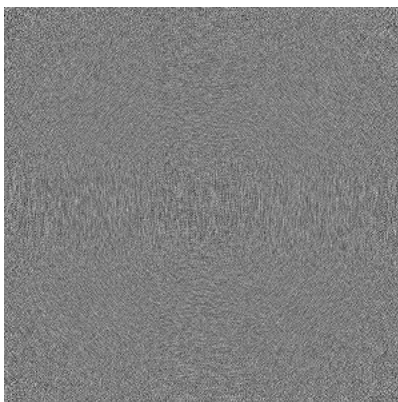


Z Index: 255

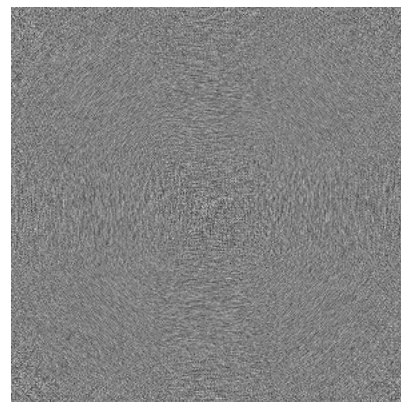
6.3.2 Raw map



X Index: 0



Y Index: 0

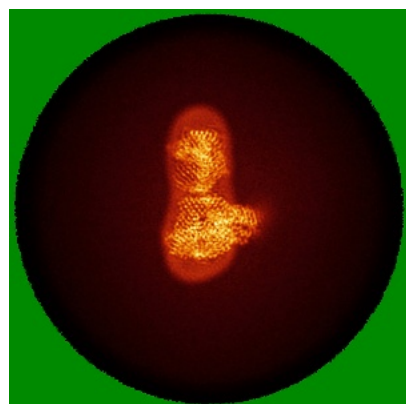


Z Index: 575

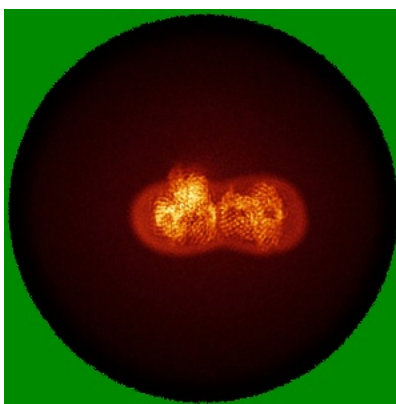
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X

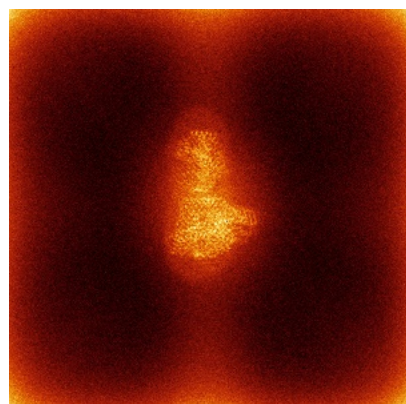


Y

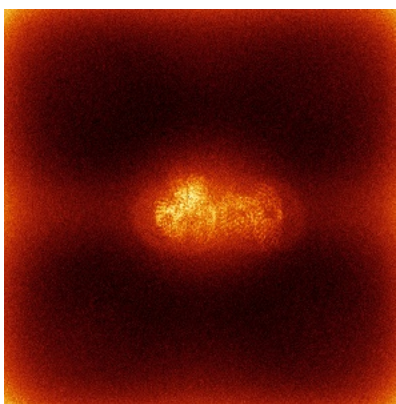


Z

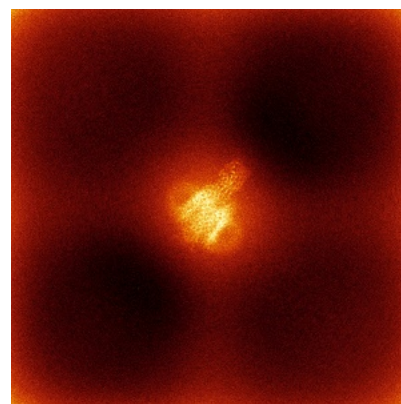
6.4.2 Raw map



X



Y

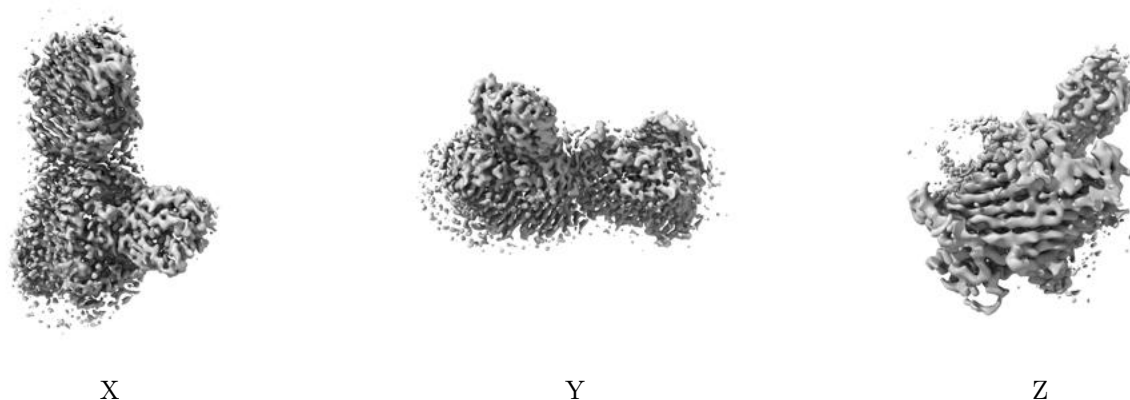


Z

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.

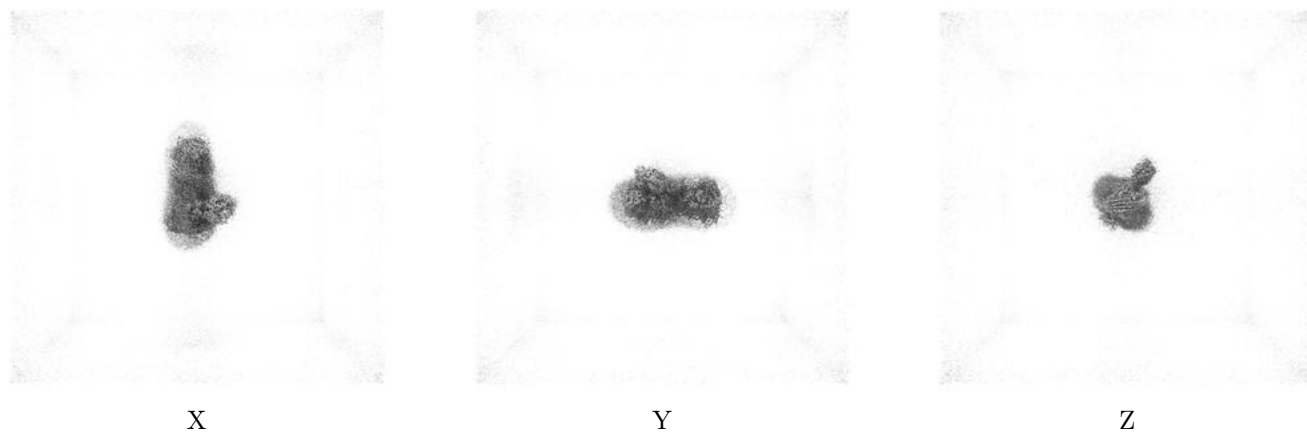
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

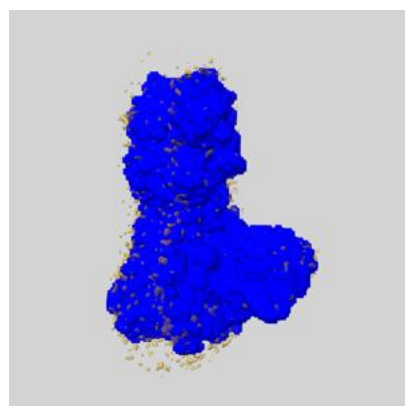
6.6 Mask visualisation [i](#)

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

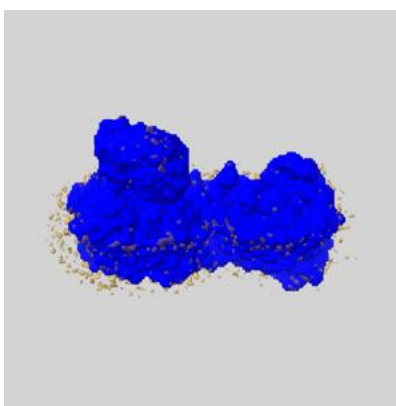
A mask typically either:

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

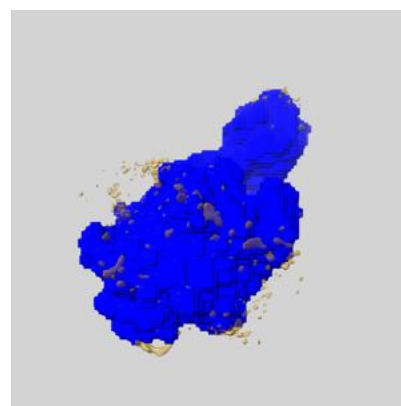
6.6.1 emd_54173_msk_1.map [i](#)



X



Y

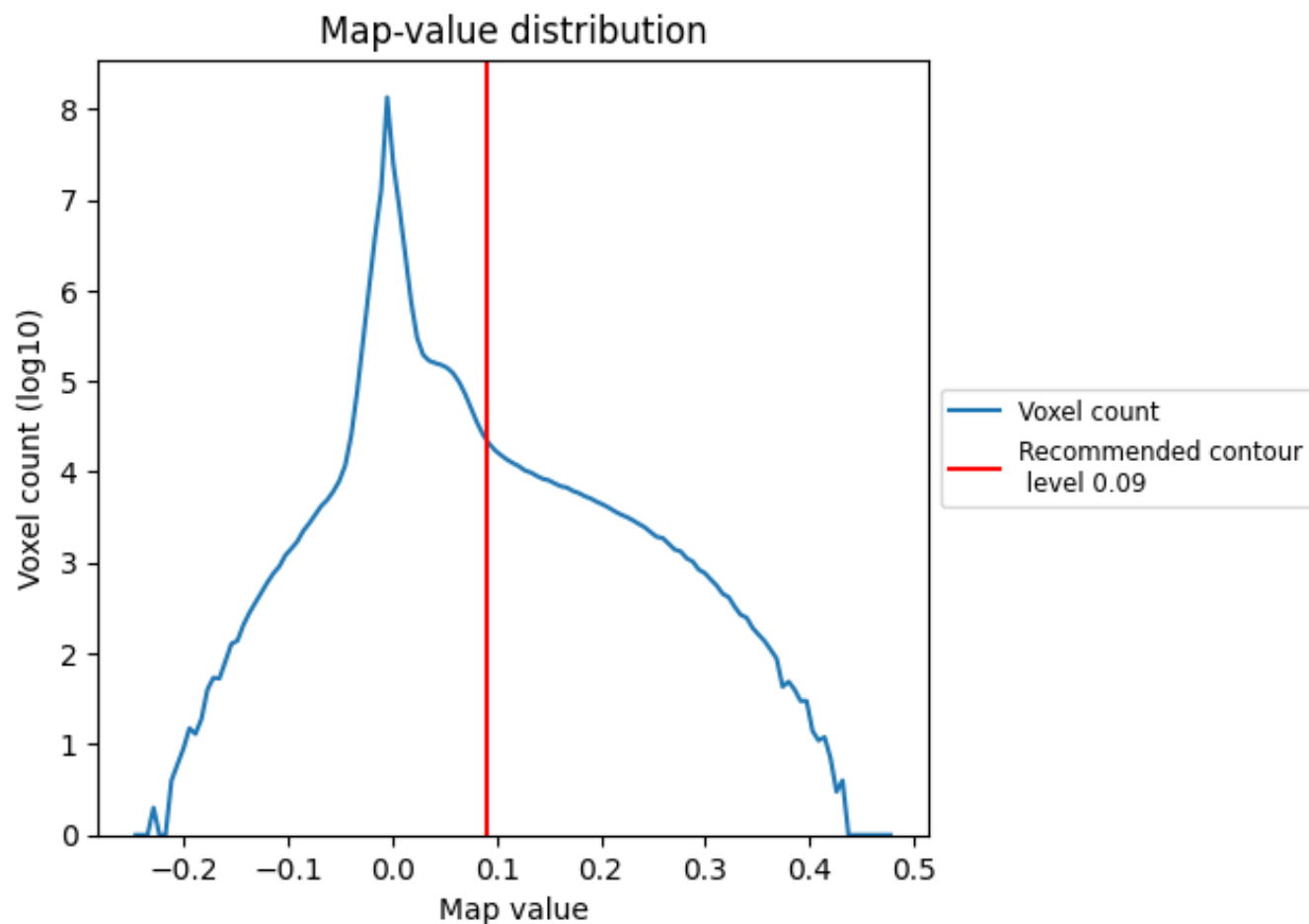


Z

7 Map analysis [i](#)

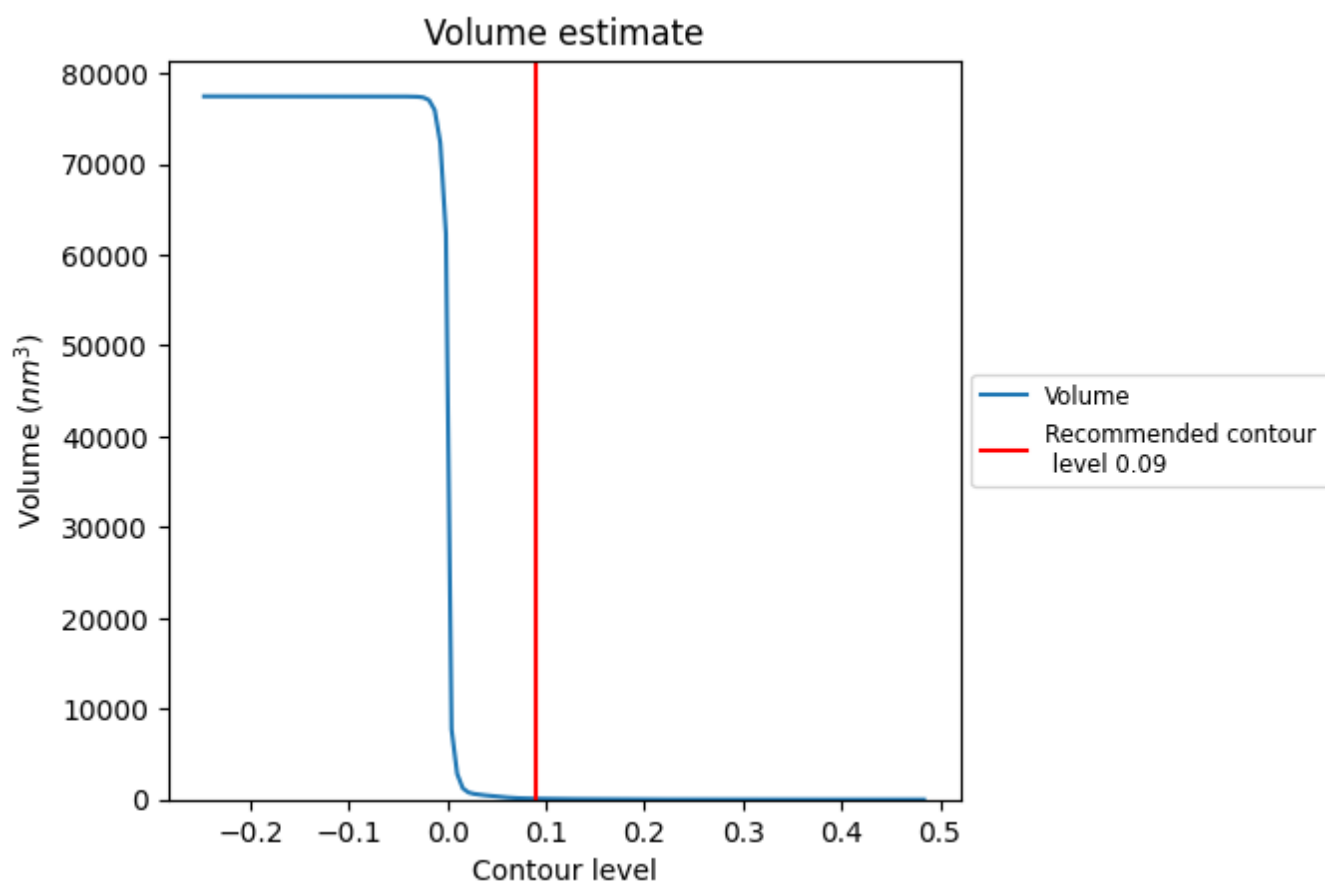
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

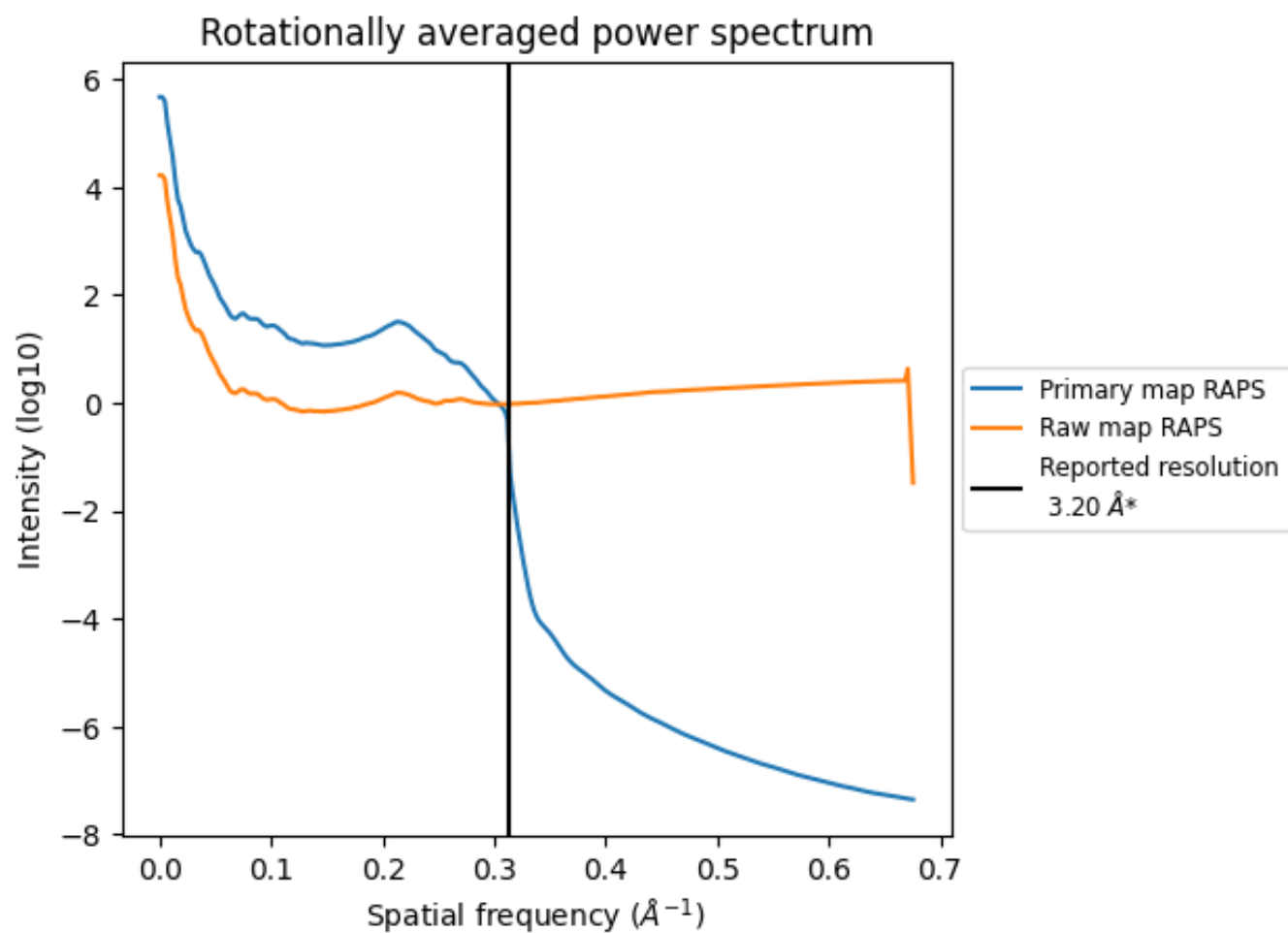
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 98 nm³; this corresponds to an approximate mass of 88 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](#)

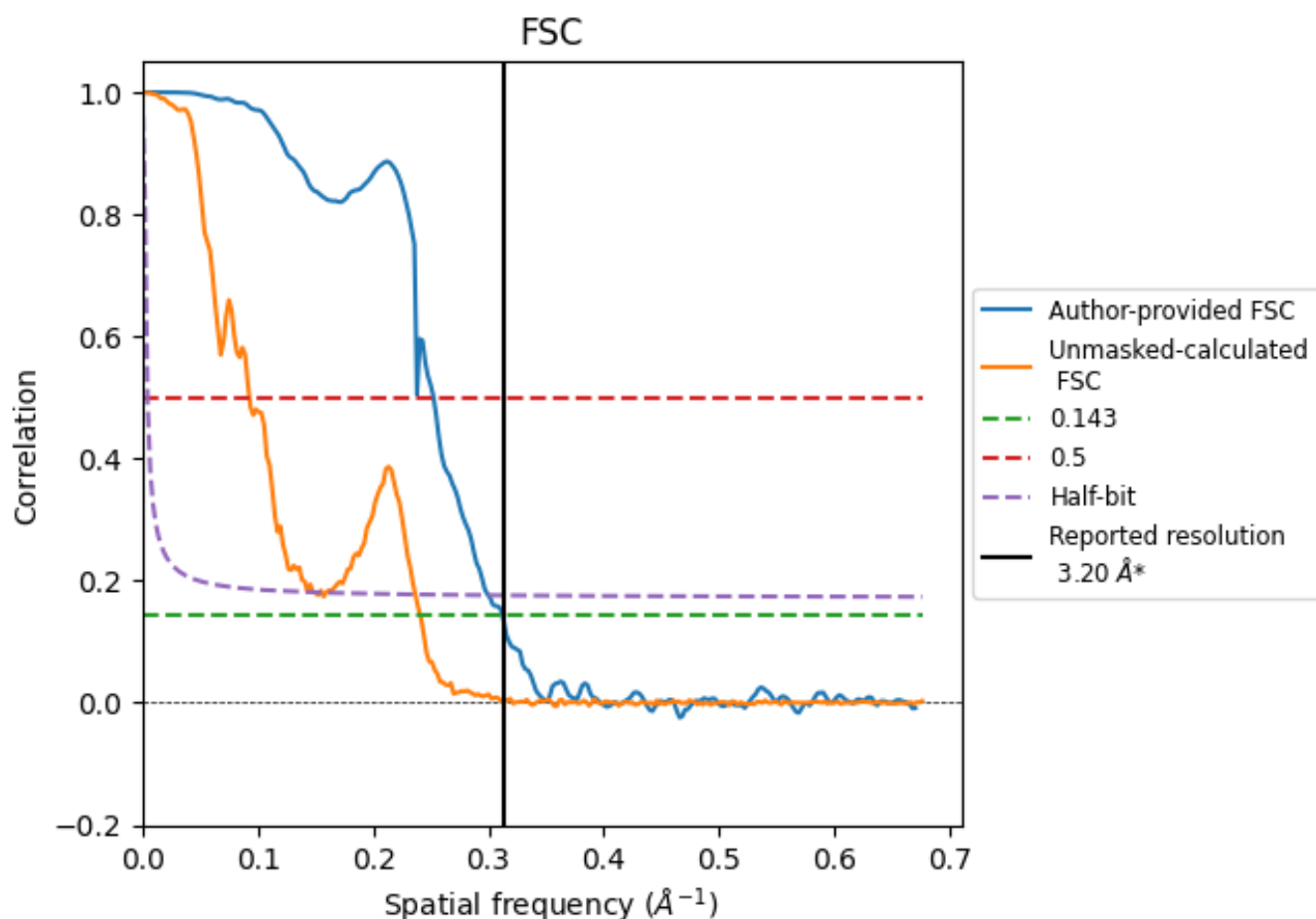


*Reported resolution corresponds to spatial frequency of 0.312 \AA^{-1}

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 0.312 \AA^{-1}

8.2 Resolution estimates [i](#)

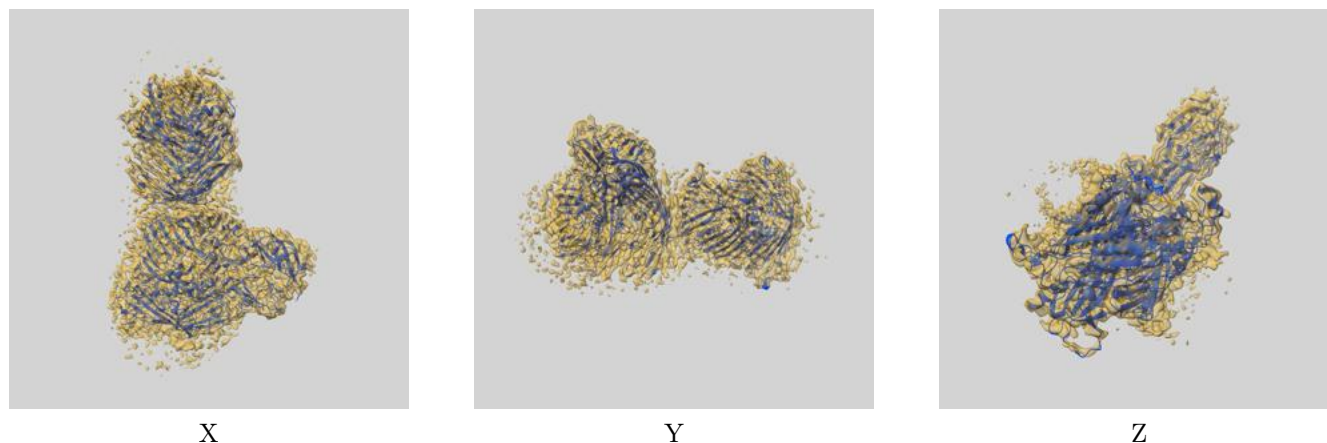
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.21	3.96	3.34
Unmasked-calculated*	4.15	10.76	6.72

*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 4.15 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-54173 and PDB model 9RPW. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



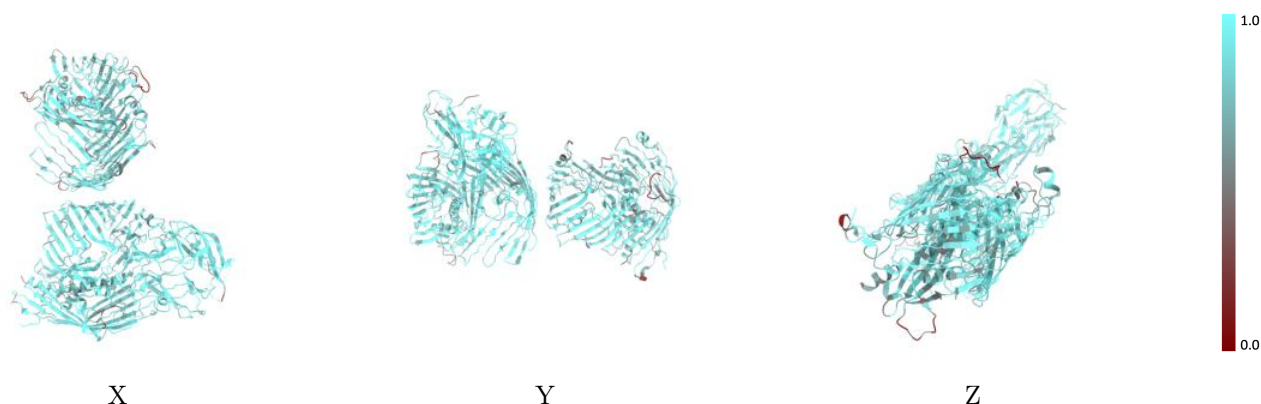
The images above show the 3D surface view of the map at the recommended contour level 0.09 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



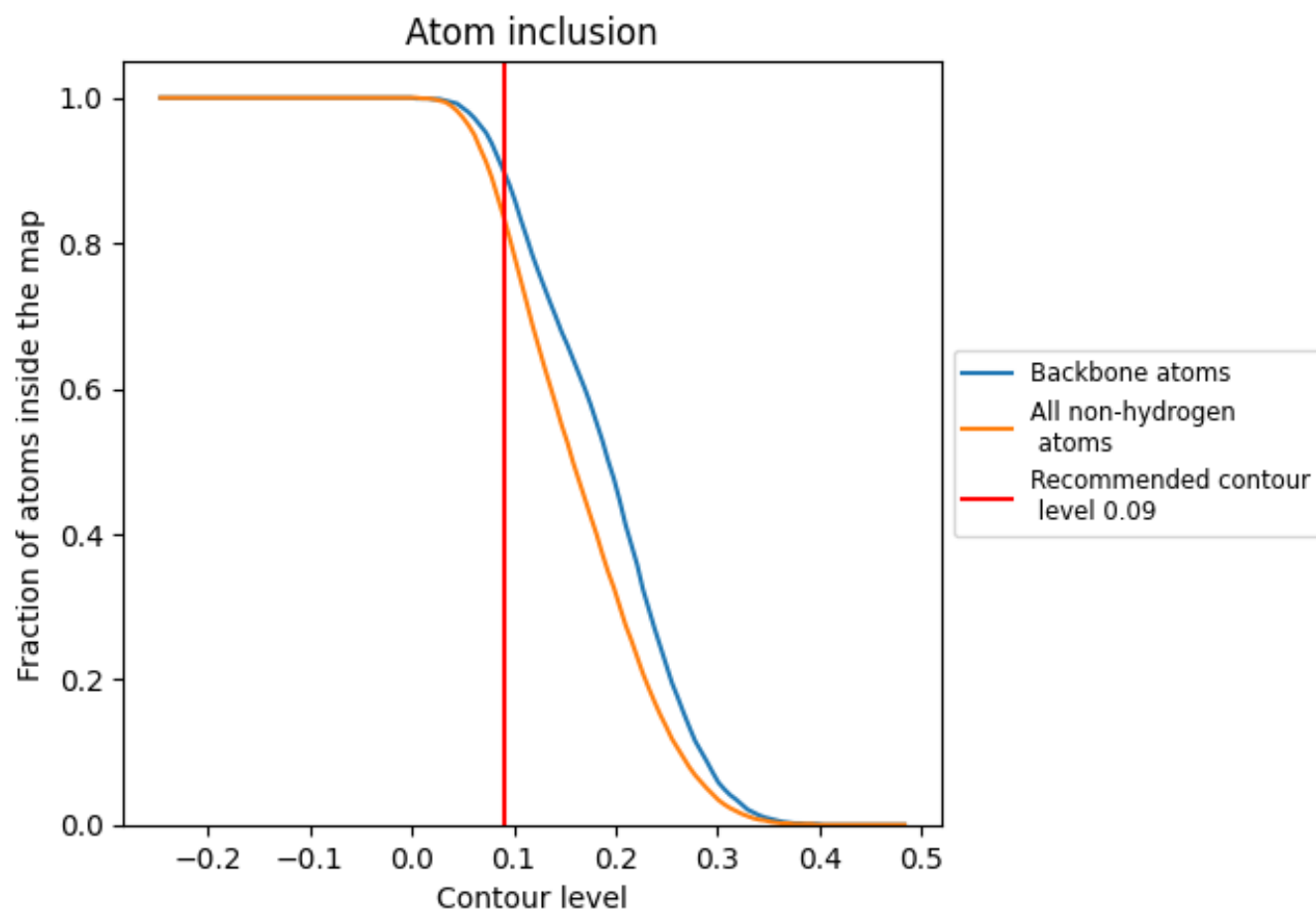
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.09).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 84% 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 (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8370	<div><div></div></div> 0.5170
A	<div><div></div></div> 0.8750	<div><div></div></div> 0.5340
B	<div><div></div></div> 0.8180	<div><div></div></div> 0.5110
C	<div><div></div></div> 0.8930	<div><div></div></div> 0.5400
D	<div><div></div></div> 0.7970	<div><div></div></div> 0.5000
E	<div><div></div></div> 0.7710	<div><div></div></div> 0.4870

