



Full wwPDB EM Validation Report ⓘ

Mar 12, 2025 – 02:13 PM JST

PDB ID : 9JUI
EMDB ID : EMD-61827
Title : Structure of Arabidopsis thaliana ABCB1 in the inward-facing conformation
Authors : Chen, Q.; Su, N.; Guo, J.
Deposited on : 2024-10-08
Resolution : 3.40 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.dev117
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.41.2

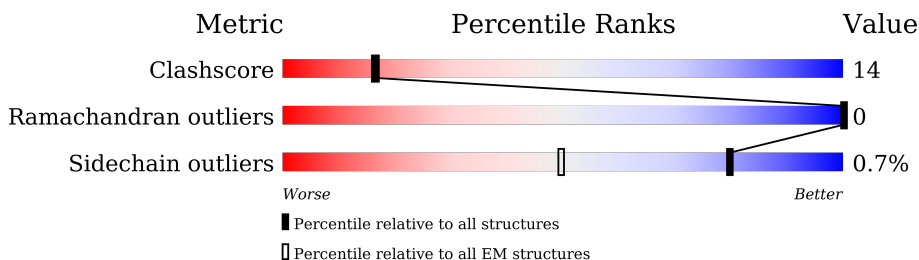
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1327	<div> <div>23%</div> <div>60%</div> <div>28%</div> <div>12%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8971 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 ABC transporter B family member 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1165	Total	C	N	O	S	0	0
			8971	5736	1543	1651	41		

There are 41 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-40	MET	-	initiating methionine	UNP Q9ZR72
A	-39	ASP	-	expression tag	UNP Q9ZR72
A	-38	TYR	-	expression tag	UNP Q9ZR72
A	-37	LYS	-	expression tag	UNP Q9ZR72
A	-36	ASP	-	expression tag	UNP Q9ZR72
A	-35	ASP	-	expression tag	UNP Q9ZR72
A	-34	ASP	-	expression tag	UNP Q9ZR72
A	-33	ASP	-	expression tag	UNP Q9ZR72
A	-32	LYS	-	expression tag	UNP Q9ZR72
A	-31	TRP	-	expression tag	UNP Q9ZR72
A	-30	SER	-	expression tag	UNP Q9ZR72
A	-29	HIS	-	expression tag	UNP Q9ZR72
A	-28	PRO	-	expression tag	UNP Q9ZR72
A	-27	GLN	-	expression tag	UNP Q9ZR72
A	-26	PHE	-	expression tag	UNP Q9ZR72
A	-25	GLU	-	expression tag	UNP Q9ZR72
A	-24	LYS	-	expression tag	UNP Q9ZR72
A	-23	GLY	-	expression tag	UNP Q9ZR72
A	-22	GLY	-	expression tag	UNP Q9ZR72
A	-21	GLY	-	expression tag	UNP Q9ZR72
A	-20	GLY	-	expression tag	UNP Q9ZR72
A	-19	SER	-	expression tag	UNP Q9ZR72
A	-18	GLY	-	expression tag	UNP Q9ZR72
A	-17	GLY	-	expression tag	UNP Q9ZR72
A	-16	SER	-	expression tag	UNP Q9ZR72
A	-15	ALA	-	expression tag	UNP Q9ZR72
A	-14	TRP	-	expression tag	UNP Q9ZR72
A	-13	SER	-	expression tag	UNP Q9ZR72

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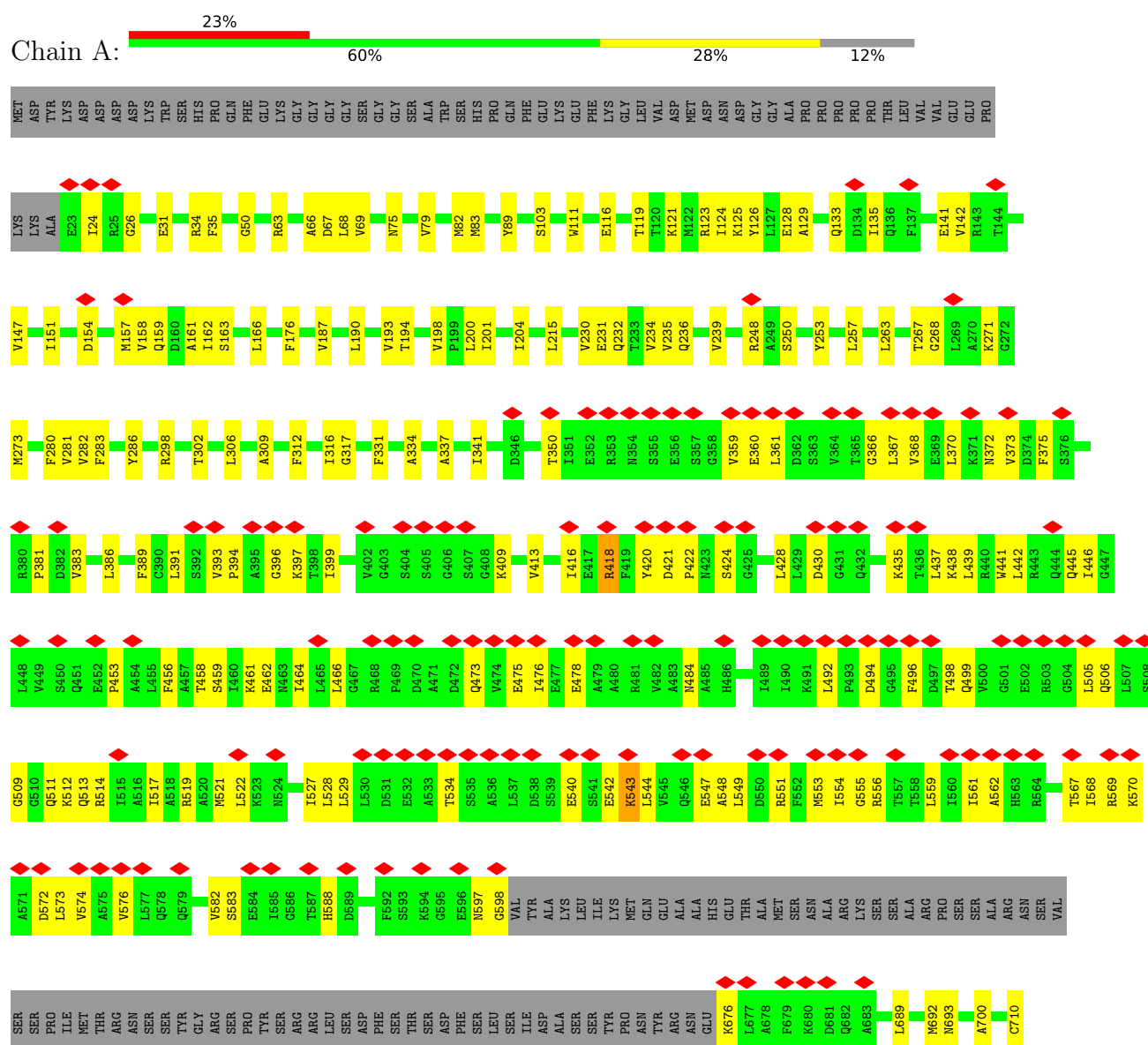
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Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	HIS	-	expression tag	UNP Q9ZR72
A	-11	PRO	-	expression tag	UNP Q9ZR72
A	-10	GLN	-	expression tag	UNP Q9ZR72
A	-9	PHE	-	expression tag	UNP Q9ZR72
A	-8	GLU	-	expression tag	UNP Q9ZR72
A	-7	LYS	-	expression tag	UNP Q9ZR72
A	-6	GLU	-	expression tag	UNP Q9ZR72
A	-5	PHE	-	expression tag	UNP Q9ZR72
A	-4	LYS	-	expression tag	UNP Q9ZR72
A	-3	GLY	-	expression tag	UNP Q9ZR72
A	-2	LEU	-	expression tag	UNP Q9ZR72
A	-1	VAL	-	expression tag	UNP Q9ZR72
A	0	ASP	-	expression tag	UNP Q9ZR72

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: ABC transporter B family member 1



A1227	H1228	V1229	I1230	A1231	V1232	I1233	D1234	D1235	G1236	K1237	V1238	A1239	E1240	Q1241	G1242	S1243	H1244	S1245	H1246	L1247	L1248	K1249	N1250	H1251	P1252	D1253	G1254	I1255	Y1256	A1257	R1258	M1259	I1260	Q1261	L1262	Q1263	R1264	PHE	THR	HIS	C209	THR	GLN	VAL	ILE	GLY	MET	THR	SER	GLY	SER	SER	SER	ARG	VAL	LYS	GLU	ASP	ALA
R1159	G1160	V1161	Q1162	L1163	S1164	G1165	G1166	Q1167	K1168	Q1169	R1170	I1171	R1175	V1178	R1179	E1182	I1183	M1184	L1185	L1186	A1189	T1190	S1191	A1192	L1193	D1194	A1195	E1196	S1197	E1198	R1199	Q1202	E1203	A1204	Q1207	A1208	C209	G1210	G1211	R1212	T1213	S1214	I1215	A1218	H1219	R1220	L1221	S1222	T1223	I1224	R1225	N1226							
L1095	K1096	A1097	I1098	R1099	K1100	H1101	I1102	A1103	I1104	E1108	P1109	C1110	L1111	F1112	G1113	T1114	T1115	I1116	Y1117	E1118	N1119	I1120	A1121	Y1122	G1123	H1124	E1125	C1126	A1127	T1128	E1129	A1130	E1131	I1132	I1133	Q1134	L1138	A1139	S1140	A1141	H1142	K1143	F1144	I1145	S1146	A1147	L1148	P1149	E1150	G1151	Y1152	K1153	T1154	Y1155	V1156	G1157	E1158		
F1031	S1032	S1035	P1037	D1038	I1039	Q1040	I1041	F1042	R1043	D1044	L1045	S1046	L1047	R1048	A1049	R1050	A1051	G1052	K1053	T1054	L1055	A1056	L1057	V1058	S1061	G1062	C1063	G1064	K1065	S1066	I1069	I1072	Q1073	R1074	F1075	Y1076	E1077	P1078	S1079	S1080	G1081	R1082	V1083	M1084	I1085	D1086	G1087	K1088	D1089	I1090	R1091	K1092	Y1093	N1094					
Y908	T909	A910	N911	L912	P915	L916	L945	A948	S949	K953	T962	V965	T979	L982	A983	R994	S995	E998	L999	L1000	K1003	T1004	E1005	I1006	E1007	P1008	D1009	D1010	P1011	D1012	T1013	T1014	P1015	V1016	P1017	D1018	R1019	L1020	R1021	G1022	E1023	V1024	E1025	L1026	K1027	H1028	I1029	D1030											
L713	F717	A723	V727	Y728	Y729	E734	Y735	M736	I737	K738	L746	L747	L750	S751	A754	L755	N758	T759	L760	Q761	H762	M765	D766	I767	V768	G769	E770	N771	L772	T773	K774	R775	V776	K779	M780	L781	V784	E788	M789	F792	E795	E796	N797	E798															
R801	I802	A807	L808	D809	A810	N811	N812	V813	R814	I817	G818	D819	R820	V823	I824	V825	Q826	G838	Q842	N843	R844	L847	V850	A851	V852	F853	P854	V855	V856	V857	D873	T882	Q883	L884	A888	R893	A896	A897	F898	N899	K903	I904	V905	R906	L907														

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	96978	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 ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1400	Depositor
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.107	Depositor
Minimum map value	-0.060	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	223.2, 223.2, 223.2	wwPDB
Map dimensions	240, 240, 240	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	0.32	0/9137	0.51	0/12366

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	8971	0	9116	252	0
All	All	8971	0	9116	252	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 14.

All (252) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:MET:HE2	1:A:337:ALA:HB1	1.50	0.93
1:A:239:VAL:HG11	1:A:1110:CYS:SG	2.25	0.77
1:A:204:ILE:HG13	1:A:281:VAL:HG21	1.67	0.76
1:A:151:ILE:HG22	1:A:882:THR:HG23	1.69	0.74
1:A:820:ARG:O	1:A:824:ILE:HG12	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:LEU:HB3	1:A:774:LYS:HD3	1.67	0.73
1:A:116:GLU:OE1	1:A:159:GLN:NE2	2.22	0.73
1:A:1104:ILE:HG22	1:A:1185:LEU:HB2	1.71	0.73
1:A:157:MET:SD	1:A:341:ILE:HD12	2.29	0.72
1:A:458:THR:HG22	1:A:459:SER:H	1.55	0.71
1:A:75:ASN:HB2	1:A:82:MET:HB3	1.72	0.70
1:A:441:TRP:O	1:A:445:GLN:NE2	2.25	0.70
1:A:373:VAL:HG11	1:A:389:PHE:HB3	1.74	0.69
1:A:842:GLN:NE2	1:A:843:TRP:O	2.26	0.68
1:A:123:ARG:NH2	1:A:151:ILE:O	2.27	0.68
1:A:234:VAL:HG12	1:A:792:PHE:HE2	1.58	0.68
1:A:372:ASN:HD22	1:A:424:SER:HB2	1.60	0.66
1:A:464:ILE:HG22	1:A:522:LEU:HD12	1.78	0.66
1:A:124:ILE:HD12	1:A:909:THR:HG22	1.78	0.65
1:A:540:GLU:HB2	1:A:543:LYS:HE3	1.78	0.65
1:A:1141:ALA:HA	1:A:1170:ARG:HD2	1.78	0.65
1:A:576:VAL:HB	1:A:583:SER:HB2	1.79	0.65
1:A:1036:ARG:HE	1:A:1039:ILE:HD13	1.61	0.65
1:A:1050:ARG:HB3	1:A:1053:LYS:HB3	1.78	0.64
1:A:1167:GLN:HG2	1:A:1170:ARG:HH21	1.61	0.63
1:A:528:LEU:HD21	1:A:549:LEU:HD12	1.80	0.63
1:A:492:LEU:HD11	1:A:505:LEU:HD21	1.80	0.63
1:A:788:GLU:OE2	1:A:1005:GLU:N	2.32	0.63
1:A:273:MET:HA	1:A:759:THR:HG22	1.80	0.62
1:A:1190:THR:HG1	1:A:1219:HIS:HE2	1.48	0.62
1:A:1115:THR:HB	1:A:1153:LYS:HB2	1.82	0.62
1:A:439:LEU:H	1:A:442:LEU:HD13	1.63	0.62
1:A:824:ILE:HD12	1:A:979:THR:HG23	1.82	0.61
1:A:1114:THR:N	1:A:1118:GLU:OE2	2.32	0.61
1:A:413:VAL:HG22	1:A:529:LEU:HD12	1.81	0.61
1:A:24:ILE:HG22	1:A:26:GLY:H	1.67	0.60
1:A:1078:PRO:HG2	1:A:1091:ARG:HH21	1.66	0.60
1:A:812:ASN:OD1	1:A:813:VAL:N	2.36	0.59
1:A:1212:ARG:HG3	1:A:1213:THR:H	1.68	0.59
1:A:775:ARG:O	1:A:779:LYS:HG2	2.03	0.59
1:A:555:GLY:O	1:A:556:ARG:HD3	2.03	0.58
1:A:809:ASP:OD1	1:A:810:ALA:N	2.37	0.58
1:A:1141:ALA:HB1	1:A:1170:ARG:HH11	1.69	0.58
1:A:119:THR:O	1:A:123:ARG:HG3	2.04	0.57
1:A:823:VAL:HG11	1:A:982:LEU:HD12	1.85	0.57
1:A:807:ALA:O	1:A:811:ASN:ND2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:LEU:HD13	1:A:391:LEU:HB2	1.88	0.56
1:A:473:GLN:HA	1:A:476:ILE:HD12	1.87	0.56
1:A:568:ILE:HG22	1:A:588:HIS:NE2	2.20	0.56
1:A:437:LEU:HB3	1:A:442:LEU:HD11	1.87	0.56
1:A:1021:ARG:CZ	1:A:1023:GLU:HB3	2.36	0.56
1:A:1220:ARG:NH2	1:A:1223:THR:O	2.38	0.55
1:A:35:PHE:HD2	1:A:121:LYS:HG3	1.71	0.55
1:A:511:GLN:HA	1:A:514:ARG:HE	1.72	0.55
1:A:394:PRO:HG2	1:A:397:LYS:HB2	1.89	0.55
1:A:1094:ASN:OD1	1:A:1095:LEU:N	2.39	0.55
1:A:1198:GLU:OE2	1:A:1202:GLN:NE2	2.39	0.55
1:A:83:MET:SD	1:A:949:SER:OG	2.57	0.55
1:A:492:LEU:HD23	1:A:494:ASP:H	1.72	0.55
1:A:567:THR:HA	1:A:570:LYS:HD3	1.88	0.55
1:A:1115:THR:N	1:A:1118:GLU:OE2	2.33	0.55
1:A:458:THR:HG22	1:A:459:SER:N	2.21	0.55
1:A:126:TYR:HD1	1:A:341:ILE:HG23	1.71	0.54
1:A:700:ALA:HB2	1:A:768:VAL:HG21	1.88	0.54
1:A:838:GLY:HA2	1:A:965:VAL:HG22	1.89	0.54
1:A:780:MET:HG2	1:A:1000:LEU:HD11	1.89	0.54
1:A:717:PHE:HD1	1:A:747:LEU:HD11	1.72	0.54
1:A:446:ILE:HG23	1:A:529:LEU:HD23	1.90	0.54
1:A:798:GLU:HB2	1:A:801:ARG:HG2	1.89	0.54
1:A:1220:ARG:HH21	1:A:1224:ILE:HA	1.73	0.54
1:A:466:LEU:HD21	1:A:519:ARG:HH11	1.73	0.53
1:A:727:VAL:HG12	1:A:736:MET:HG3	1.89	0.53
1:A:693:ASN:HB2	1:A:772:LEU:HD13	1.90	0.53
1:A:1102:ILE:HG13	1:A:1183:ILE:HB	1.89	0.53
1:A:232:GLN:O	1:A:236:GLN:HG2	2.09	0.53
1:A:1167:GLN:O	1:A:1170:ARG:HG2	2.08	0.53
1:A:1186:LEU:N	1:A:1215:ILE:O	2.38	0.53
1:A:268:GLY:HA2	1:A:766:ASP:OD2	2.09	0.53
1:A:1021:ARG:CZ	1:A:1086:ASP:HA	2.39	0.53
1:A:461:LYS:NZ	1:A:496:PHE:O	2.36	0.53
1:A:574:VAL:HB	1:A:588:HIS:HD2	1.73	0.53
1:A:126:TYR:CD1	1:A:341:ILE:HG23	2.43	0.52
1:A:360:GLU:HA	1:A:441:TRP:CE3	2.45	0.52
1:A:761:GLN:HG3	1:A:762:HIS:N	2.25	0.52
1:A:819:ASP:OD2	1:A:820:ARG:NH1	2.42	0.52
1:A:453:PRO:HG2	1:A:512:LYS:HE2	1.91	0.52
1:A:797:ASN:HB3	1:A:802:ILE:HD11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:905:VAL:O	1:A:909:THR:HG23	2.09	0.52
1:A:854:PRO:HA	1:A:857:VAL:HG22	1.91	0.52
1:A:1220:ARG:HB3	1:A:1224:ILE:HD12	1.91	0.52
1:A:413:VAL:O	1:A:416:ILE:HG22	2.09	0.52
1:A:994:ARG:HD2	1:A:995:SER:N	2.25	0.52
1:A:521:MET:SD	1:A:549:LEU:HD21	2.49	0.51
1:A:215:LEU:HB3	1:A:267:THR:HG22	1.92	0.51
1:A:776:VAL:O	1:A:780:MET:HG3	2.11	0.51
1:A:194:THR:O	1:A:198:VAL:HG23	2.10	0.51
1:A:31:GLU:O	1:A:34:ARG:HG2	2.10	0.51
1:A:1021:ARG:H	1:A:1021:ARG:HD3	1.76	0.51
1:A:190:LEU:HA	1:A:193:VAL:HG22	1.93	0.51
1:A:754:ALA:O	1:A:758:ASN:HB2	2.11	0.51
1:A:713:LEU:HD23	1:A:750:LEU:HD22	1.91	0.51
1:A:1057:LEU:HA	1:A:1230:ILE:HG23	1.93	0.50
1:A:1139:ALA:HA	1:A:1204:ALA:HB1	1.93	0.50
1:A:713:LEU:HD13	1:A:754:ALA:HA	1.93	0.50
1:A:124:ILE:HD13	1:A:908:TYR:CD2	2.47	0.50
1:A:201:ILE:HD13	1:A:282:VAL:HG22	1.94	0.49
1:A:373:VAL:HG13	1:A:386:LEU:HB3	1.93	0.49
1:A:253:TYR:O	1:A:257:LEU:HG	2.13	0.49
1:A:506:GLN:HG3	1:A:509:GLY:H	1.78	0.49
1:A:844:ARG:HA	1:A:847:LEU:HB2	1.94	0.49
1:A:1038:ASP:OD1	1:A:1039:ILE:HD12	2.12	0.49
1:A:1199:ARG:HG2	1:A:1199:ARG:HH11	1.77	0.49
1:A:770:GLU:O	1:A:774:LYS:HG2	2.12	0.49
1:A:542:GLU:OE2	1:A:542:GLU:N	2.28	0.49
1:A:850:VAL:HA	1:A:853:PHE:HB2	1.95	0.49
1:A:1165:GLY:O	1:A:1169:GLN:HG2	2.13	0.49
1:A:1148:LEU:HD13	1:A:1154:THR:HG23	1.94	0.48
1:A:572:ASP:OD1	1:A:573:LEU:N	2.47	0.48
1:A:66:ALA:HB2	1:A:309:ALA:HB2	1.96	0.48
1:A:250:SER:HA	1:A:781:LEU:HD23	1.96	0.48
1:A:466:LEU:HD21	1:A:519:ARG:NH1	2.27	0.48
1:A:476:ILE:HA	1:A:522:LEU:HD11	1.95	0.48
1:A:125:LYS:O	1:A:128:GLU:HG3	2.14	0.48
1:A:1043:ARG:N	1:A:1236:GLY:O	2.47	0.48
1:A:1117:TYR:CE1	1:A:1132:ILE:HD12	2.48	0.48
1:A:758:ASN:HA	1:A:761:GLN:HG2	1.96	0.48
1:A:1103:ALA:HB2	1:A:1179:ARG:HG2	1.95	0.48
1:A:157:MET:O	1:A:334:ALA:HB1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1027:LYS:HD3	1:A:1028:HIS:HB2	1.95	0.47
1:A:1095:LEU:O	1:A:1099:ARG:HB2	2.14	0.47
1:A:747:LEU:HD12	1:A:750:LEU:HD12	1.95	0.47
1:A:283:PHE:HB2	1:A:751:SER:HB3	1.96	0.47
1:A:801:ARG:HE	1:A:801:ARG:HA	1.79	0.47
1:A:66:ALA:HB1	1:A:306:LEU:HD23	1.97	0.47
1:A:1256:TYR:CE2	1:A:1260:ILE:HD11	2.49	0.47
1:A:727:VAL:O	1:A:736:MET:HG3	2.14	0.47
1:A:907:LEU:O	1:A:911:ASN:ND2	2.47	0.47
1:A:540:GLU:O	1:A:543:LYS:HG3	2.14	0.47
1:A:476:ILE:HG12	1:A:522:LEU:HD11	1.96	0.47
1:A:1016:VAL:HG21	1:A:1098:ILE:HG12	1.97	0.47
1:A:389:PHE:HD1	1:A:582:VAL:HB	1.80	0.46
1:A:689:LEU:HD21	1:A:772:LEU:HD21	1.96	0.46
1:A:781:LEU:HA	1:A:784:VAL:HG12	1.96	0.46
1:A:1074:ARG:HH11	1:A:1091:ARG:HA	1.79	0.46
1:A:422:PRO:HG3	1:A:435:LYS:HE2	1.97	0.46
1:A:547:GLU:O	1:A:551:ARG:NH2	2.32	0.46
1:A:79:VAL:HB	1:A:953:LYS:HG2	1.96	0.46
1:A:129:ALA:O	1:A:133:GLN:HG2	2.14	0.46
1:A:393:VAL:HG12	1:A:399:ILE:HD13	1.97	0.46
1:A:734:GLU:O	1:A:738:LYS:HG2	2.15	0.46
1:A:1175:ARG:HA	1:A:1178:VAL:HG12	1.98	0.46
1:A:1080:SER:O	1:A:1082:ARG:NH1	2.49	0.46
1:A:409:LYS:HE2	1:A:561:ILE:HD11	1.97	0.46
1:A:903:LYS:NZ	1:A:906:ARG:HE	2.13	0.46
1:A:544:LEU:O	1:A:548:ALA:N	2.49	0.46
1:A:798:GLU:O	1:A:802:ILE:HG12	2.16	0.46
1:A:912:LEU:C	1:A:915:PRO:HD2	2.36	0.46
1:A:758:ASN:OD1	1:A:761:GLN:NE2	2.49	0.45
1:A:823:VAL:HA	1:A:826:GLN:HG2	1.98	0.45
1:A:366:GLY:HA2	1:A:430:ASP:HB3	1.98	0.45
1:A:710:CYS:HA	1:A:713:LEU:HD12	1.97	0.45
1:A:34:ARG:HG3	1:A:35:PHE:CD1	2.51	0.45
1:A:1228:HIS:ND1	1:A:1229:VAL:HG13	2.31	0.45
1:A:1120:ILE:HG22	1:A:1178:VAL:HG11	1.98	0.45
1:A:1175:ARG:O	1:A:1178:VAL:HG12	2.17	0.45
1:A:597:ASN:OD1	1:A:598:GLY:N	2.49	0.45
1:A:375:PHE:CE1	1:A:421:ASP:HB2	2.52	0.45
1:A:1109:PRO:O	1:A:1168:LYS:HD2	2.16	0.45
1:A:723:ALA:O	1:A:727:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:773:THR:HB	1:A:814:ARG:HB2	1.98	0.45
1:A:68:LEU:HD22	1:A:945:LEU:HD13	1.98	0.45
1:A:280:PHE:HA	1:A:751:SER:HB2	1.99	0.45
1:A:368:VAL:HA	1:A:428:LEU:O	2.17	0.45
1:A:553:MET:SD	1:A:554:ILE:N	2.90	0.45
1:A:475:GLU:HA	1:A:478:GLU:HB2	1.98	0.44
1:A:755:LEU:O	1:A:759:THR:HG23	2.17	0.44
1:A:418:ARG:HD2	1:A:439:LEU:HD13	1.99	0.44
1:A:574:VAL:HB	1:A:588:HIS:CD2	2.51	0.44
1:A:689:LEU:HD11	1:A:772:LEU:HD11	1.99	0.44
1:A:1131:GLU:O	1:A:1134:GLN:HG3	2.17	0.44
1:A:494:ASP:OD2	1:A:498:THR:HG22	2.18	0.44
1:A:200:LEU:O	1:A:204:ILE:HG12	2.18	0.44
1:A:534:THR:HG22	1:A:562:ALA:HB2	2.00	0.44
1:A:141:GLU:HG2	1:A:142:VAL:H	1.81	0.44
1:A:1254:GLY:O	1:A:1258:ARG:NH1	2.26	0.44
1:A:350:THR:HG22	1:A:420:TYR:HE2	1.83	0.43
1:A:1109:PRO:HG2	1:A:1168:LYS:HG2	1.99	0.43
1:A:1145:ILE:HG22	1:A:1152:TYR:HB2	2.00	0.43
1:A:1120:ILE:HB	1:A:1132:ILE:HD13	2.00	0.43
1:A:381:PRO:O	1:A:383:VAL:HG13	2.18	0.43
1:A:948:ALA:HB1	1:A:962:THR:HB	2.01	0.43
1:A:312:PHE:O	1:A:316:ILE:HG12	2.18	0.43
1:A:462:GLU:OE2	1:A:903:LYS:HD2	2.19	0.43
1:A:162:ILE:HA	1:A:166:LEU:HB2	2.00	0.43
1:A:298:ARG:HD3	1:A:737:ILE:HD11	2.00	0.43
1:A:135:ILE:HD12	1:A:893:ARG:HG2	2.00	0.43
1:A:484:ASN:H	1:A:548:ALA:HB2	1.84	0.43
1:A:1124:HIS:HD2	1:A:1126:CYS:HB2	1.82	0.43
1:A:187:VAL:HG22	1:A:302:THR:HG21	2.01	0.42
1:A:232:GLN:HA	1:A:235:VAL:HG22	2.00	0.42
1:A:372:ASN:OD1	1:A:373:VAL:N	2.51	0.42
1:A:1109:PRO:HG2	1:A:1168:LYS:HE3	2.01	0.42
1:A:1261:GLN:HA	1:A:1264:ARG:HE	1.84	0.42
1:A:994:ARG:O	1:A:998:GLU:HG2	2.20	0.42
1:A:1047:LEU:HD11	1:A:1055:LEU:HD23	2.00	0.42
1:A:1054:THR:OG1	1:A:1209:CYS:HB3	2.20	0.42
1:A:458:THR:CG2	1:A:459:SER:H	2.27	0.42
1:A:1045:LEU:HD21	1:A:1047:LEU:HB2	2.01	0.42
1:A:63:ARG:NH1	1:A:67:ASP:OD1	2.53	0.42
1:A:359:VAL:HG12	1:A:361:LEU:H	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:GLY:N	1:A:556:ARG:HH22	2.17	0.42
1:A:789:MET:HB2	1:A:1075:PHE:O	2.20	0.42
1:A:234:VAL:HG12	1:A:792:PHE:CE2	2.47	0.41
1:A:438:LYS:O	1:A:439:LEU:HB3	2.19	0.41
1:A:1199:ARG:HG2	1:A:1199:ARG:NH1	2.35	0.41
1:A:50:GLY:O	1:A:103:SER:HB3	2.20	0.41
1:A:111:TRP:HB3	1:A:163:SER:O	2.20	0.41
1:A:230:VAL:O	1:A:234:VAL:HG22	2.19	0.41
1:A:372:ASN:ND2	1:A:424:SER:HB2	2.32	0.41
1:A:126:TYR:OH	1:A:147:VAL:HG13	2.20	0.41
1:A:569:ARG:NH1	1:A:588:HIS:CE1	2.89	0.41
1:A:161:ALA:O	1:A:166:LEU:N	2.53	0.41
1:A:367:LEU:H	1:A:430:ASP:HB3	1.84	0.41
1:A:396:GLY:HA2	1:A:554:ILE:HA	2.01	0.41
1:A:689:LEU:HA	1:A:692:MET:HG2	2.02	0.41
1:A:69:VAL:HG12	1:A:729:TYR:HE2	1.85	0.41
1:A:250:SER:HA	1:A:781:LEU:CD2	2.50	0.41
1:A:1194:ASP:OD1	1:A:1195:ALA:N	2.53	0.41
1:A:271:LYS:HE3	1:A:762:HIS:ND1	2.35	0.41
1:A:459:SER:HB2	1:A:499:GLN:OE1	2.21	0.41
1:A:527:ILE:HA	1:A:556:ARG:HG3	2.03	0.41
1:A:746:LEU:O	1:A:750:LEU:HG	2.21	0.41
1:A:123:ARG:CZ	1:A:154:ASP:HB3	2.51	0.41
1:A:166:LEU:HA	1:A:331:PHE:HZ	1.85	0.41
1:A:176:PHE:HA	1:A:317:GLY:HA2	2.02	0.41
1:A:484:ASN:HD21	1:A:517:ILE:HB	1.85	0.41
1:A:852:VAL:O	1:A:856:VAL:HG23	2.21	0.41
1:A:884:LEU:HD13	1:A:911:ASN:ND2	2.35	0.41
1:A:67:ASP:HB3	1:A:89:TYR:CZ	2.56	0.41
1:A:116:GLU:HG3	1:A:916:LEU:HD12	2.02	0.41
1:A:1038:ASP:OD1	1:A:1038:ASP:N	2.53	0.41
1:A:1116:ILE:HA	1:A:1119:ASN:ND2	2.36	0.40
1:A:119:THR:OG1	1:A:158:VAL:HG13	2.21	0.40
1:A:453:PRO:HG3	1:A:513:GLN:OE1	2.21	0.40
1:A:456:PHE:HE1	1:A:888:ALA:HA	1.86	0.40
1:A:1134:GLN:O	1:A:1138:LEU:HG	2.21	0.40
1:A:1195:ALA:O	1:A:1199:ARG:HG3	2.22	0.40
1:A:236:GLN:NE2	1:A:1112:PHE:CE1	2.89	0.40
1:A:409:LYS:HD2	1:A:559:LEU:HD13	2.02	0.40
1:A:231:GLU:O	1:A:235:VAL:HG13	2.22	0.40
1:A:765:TRP:HA	1:A:768:VAL:HG12	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1052:GLY:HA3	1:A:1212:ARG:HH21	1.86	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	1161/1327 (88%)	1106 (95%)	55 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	949/1087 (87%)	942 (99%)	7 (1%)	81	88

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	248	ARG
1	A	263	LEU
1	A	286	TYR
1	A	418	ARG
1	A	543	LYS

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Mol	Chain	Res	Type
1	A	676	LYS
1	A	1021	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	HIS
1	A	236	GLN
1	A	388	ASN
1	A	445	GLN
1	A	484	ASN
1	A	1124	HIS

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.

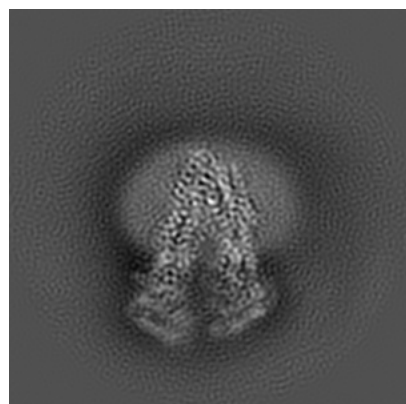
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61827. These allow visual inspection of the internal detail of the map and identification of artifacts.

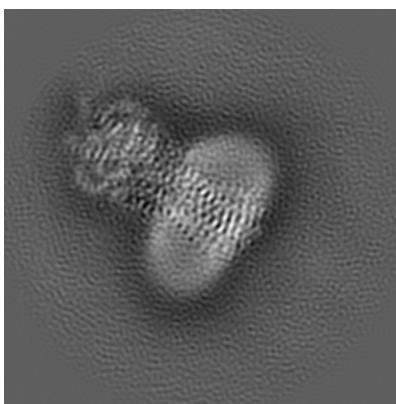
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

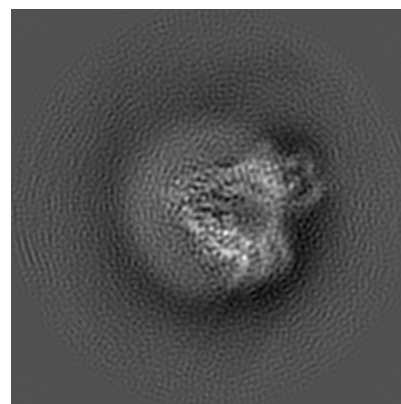
6.1.1 Primary map



X

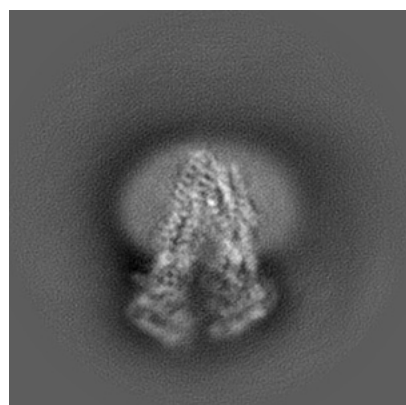


Y

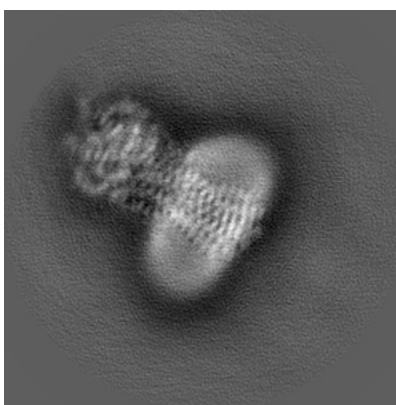


Z

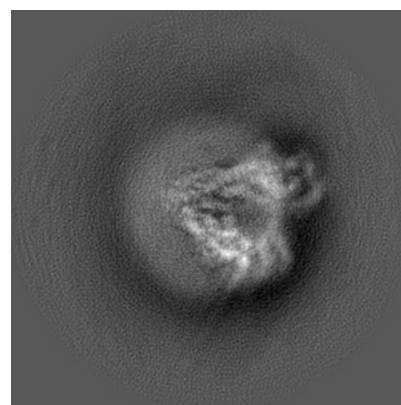
6.1.2 Raw map



X



Y

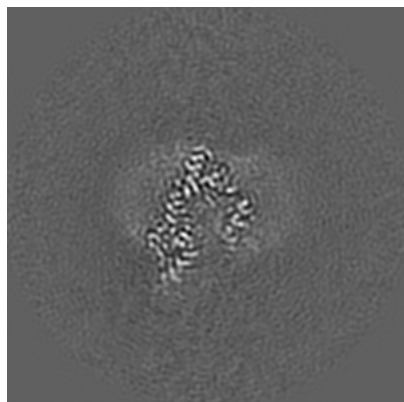


Z

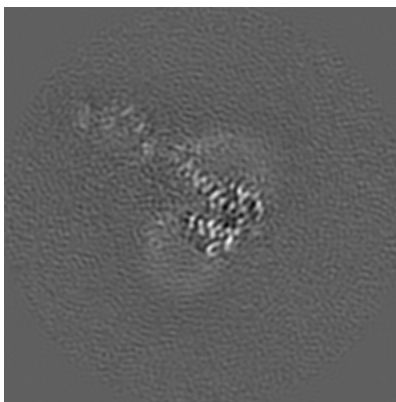
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

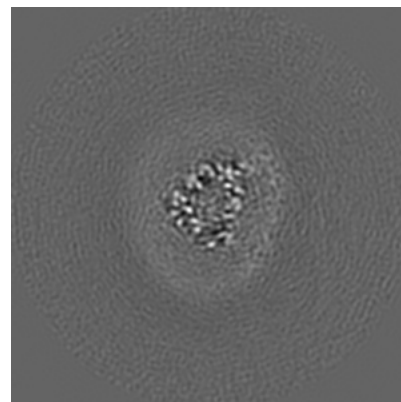
6.2.1 Primary map



X Index: 120

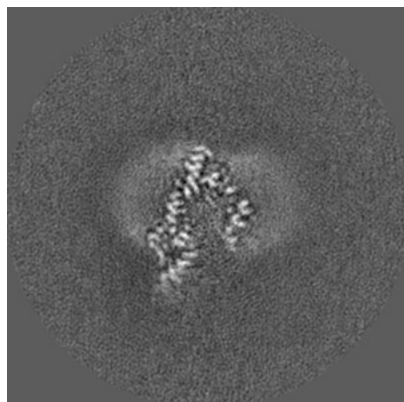


Y Index: 120



Z Index: 120

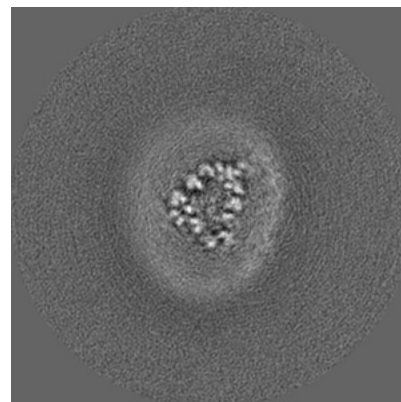
6.2.2 Raw map



X Index: 120



Y Index: 120

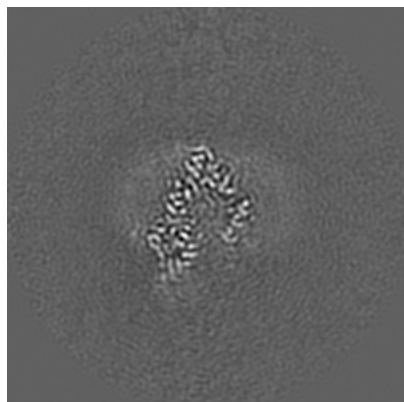


Z Index: 120

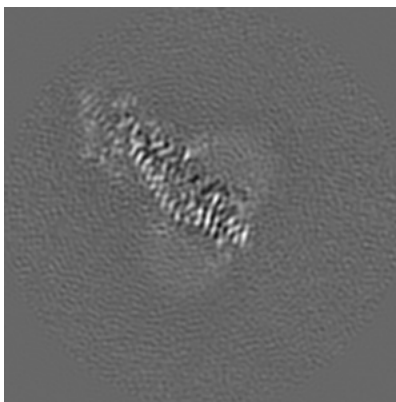
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

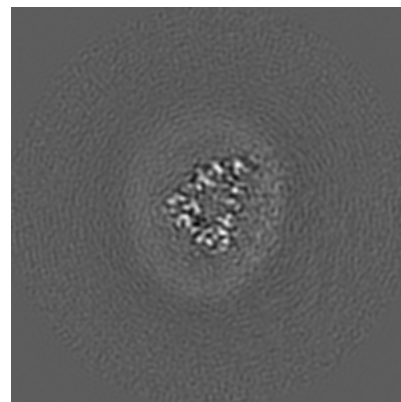
6.3.1 Primary map



X Index: 119

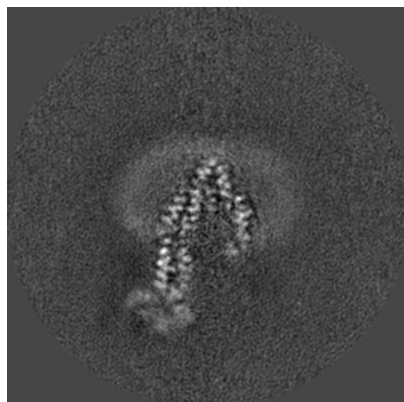


Y Index: 135

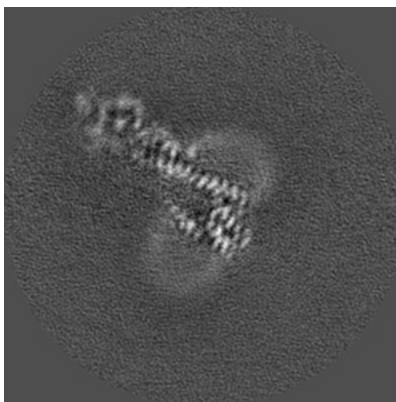


Z Index: 118

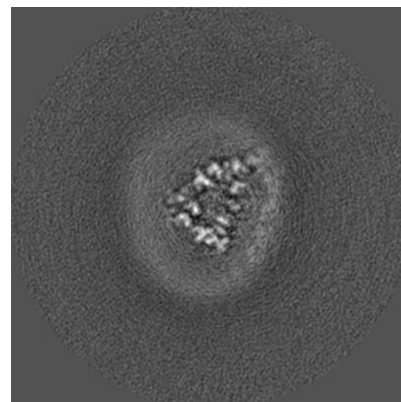
6.3.2 Raw map



X Index: 129



Y Index: 130

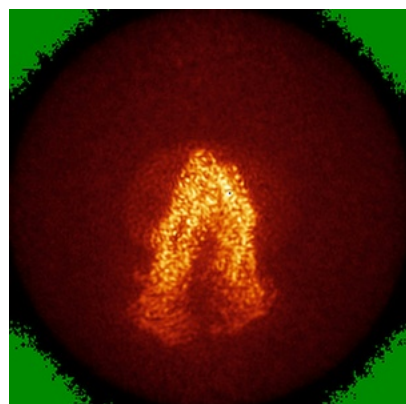


Z Index: 117

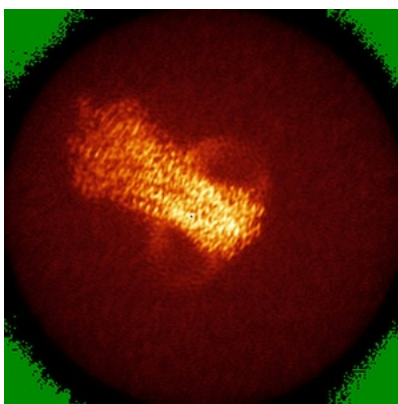
The images above show the largest variance slices of the map in three orthogonal directions.

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

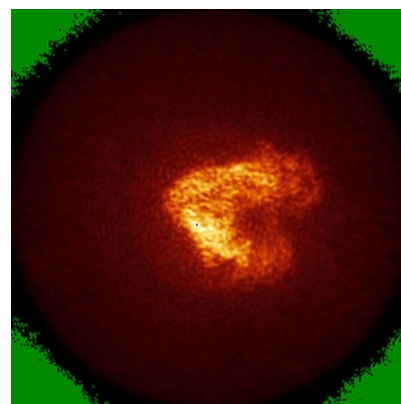
6.4.1 Primary map



X

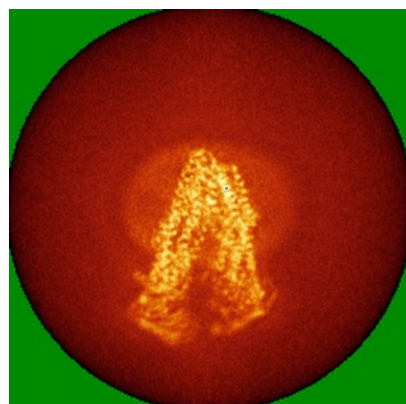


Y

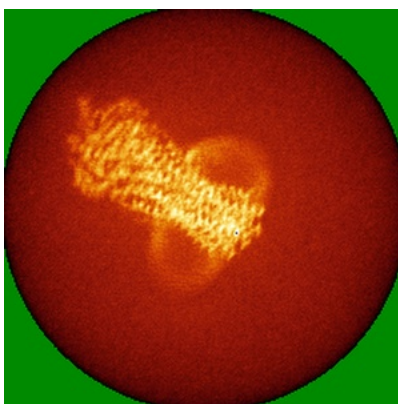


Z

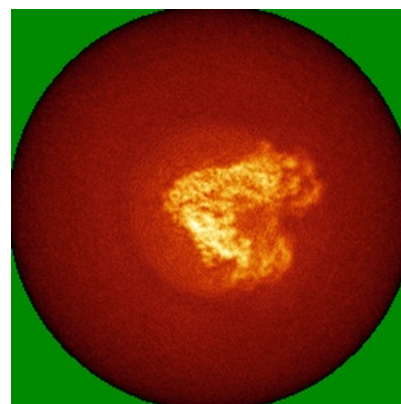
6.4.2 Raw map



X



Y

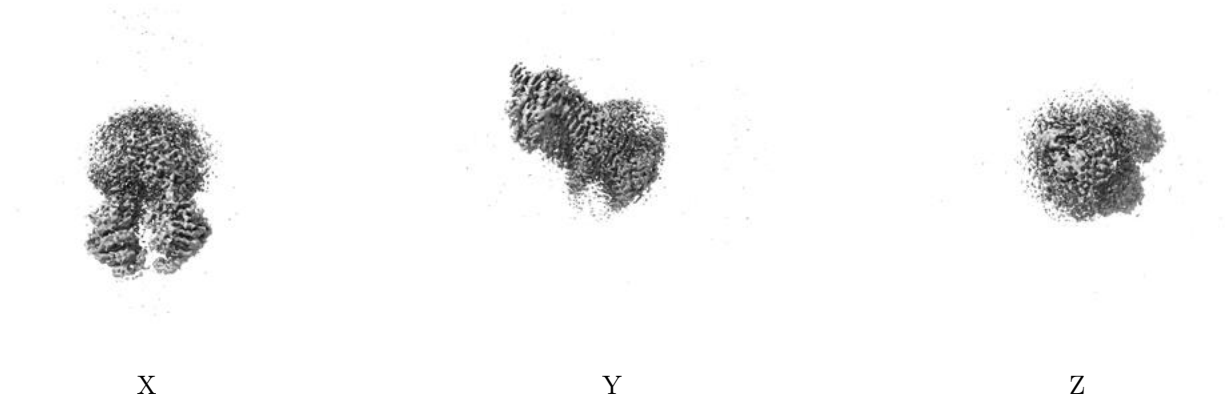


Z

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.

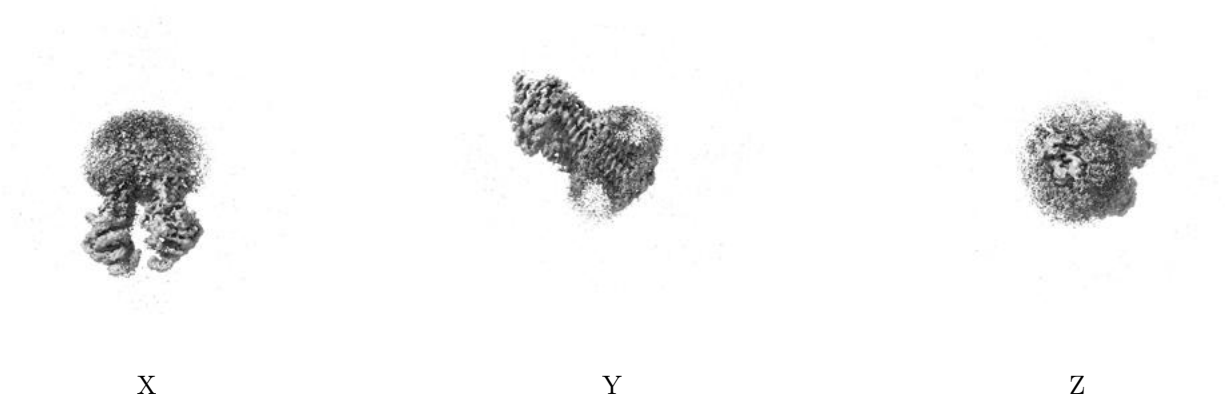
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.01. 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.

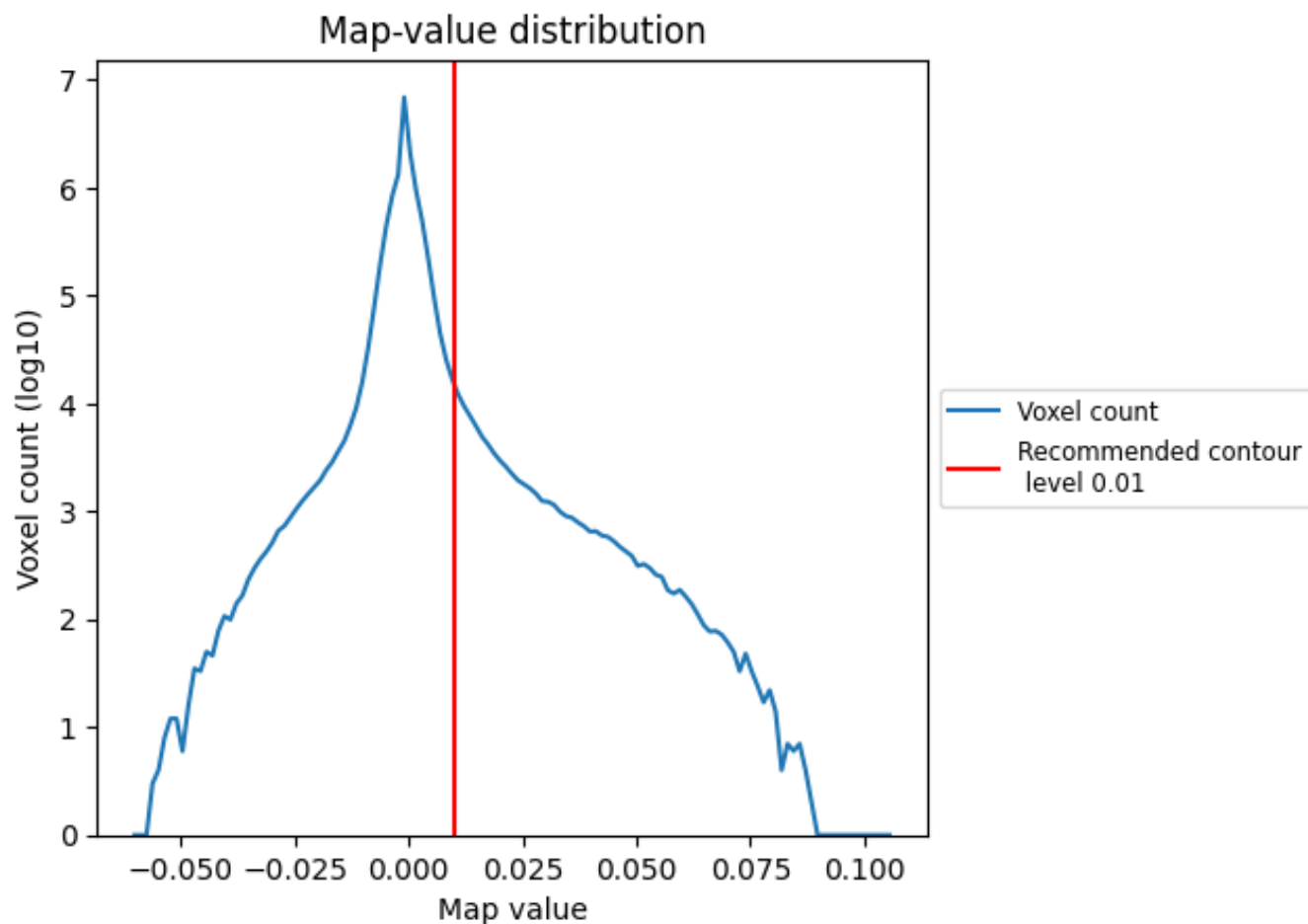
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

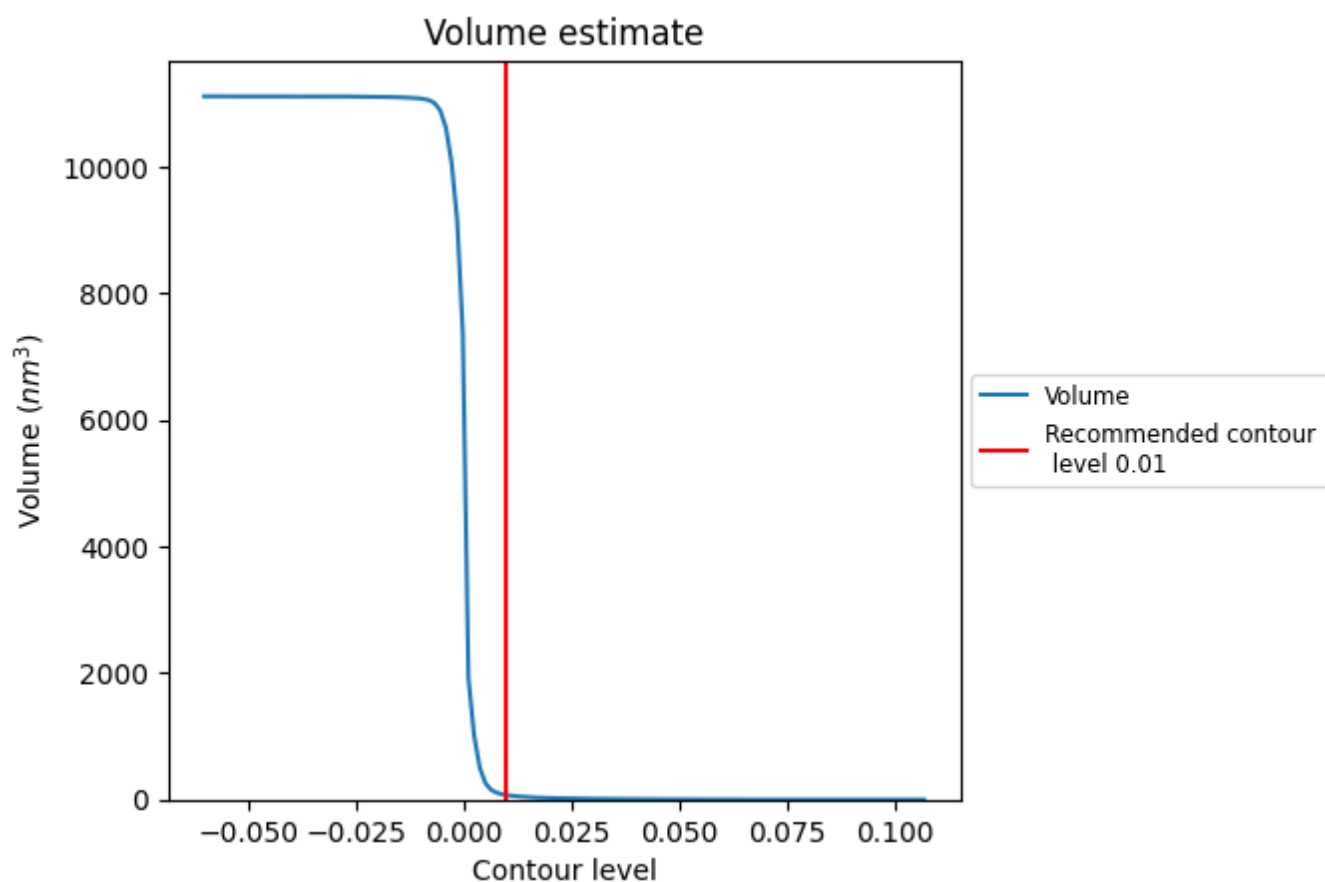
This section contains the results of statistical analysis of the map.

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.

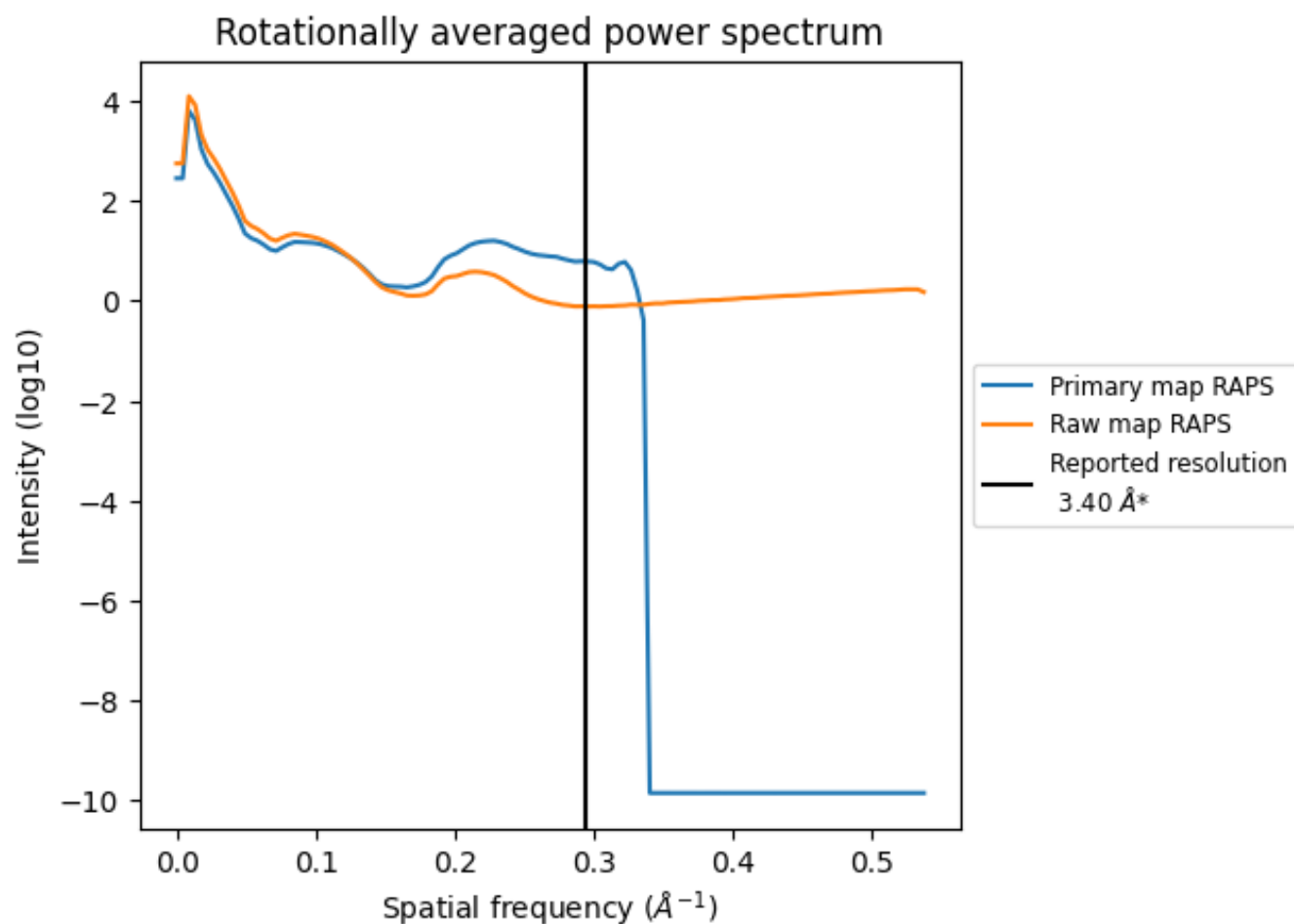
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 71 nm^3 ; this corresponds to an approximate mass of 64 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 ⓘ

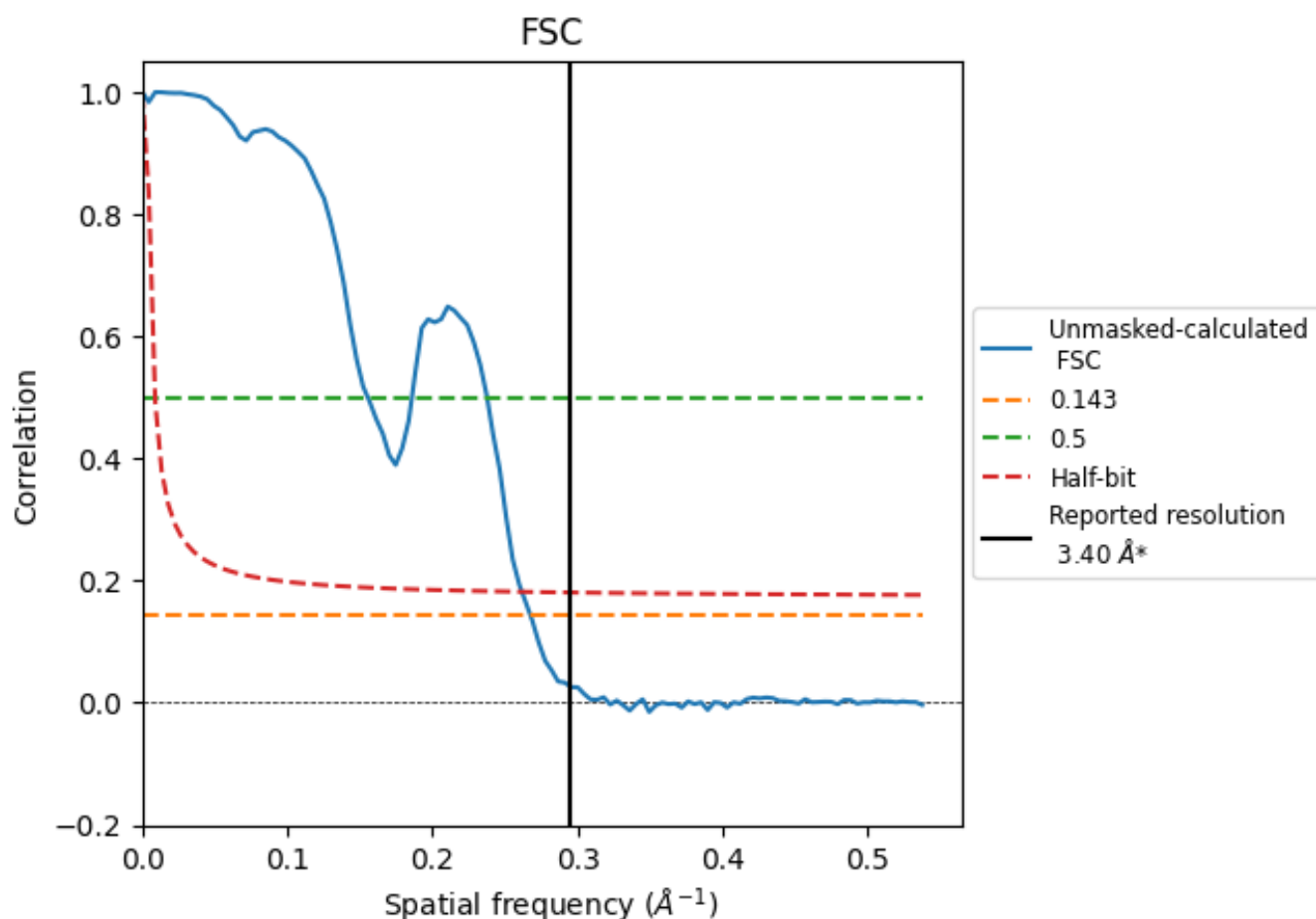


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

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 0.294 \AA^{-1}

8.2 Resolution estimates [i](#)

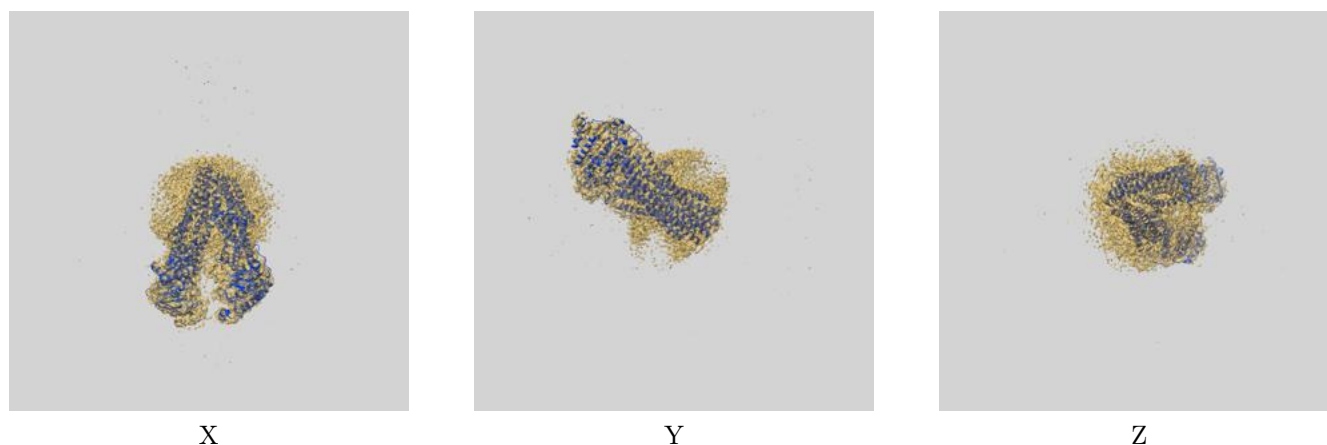
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.74	6.43	3.82

*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 3.74 differs from the reported value 3.4 by more than 10 %

9 Map-model fit [i](#)

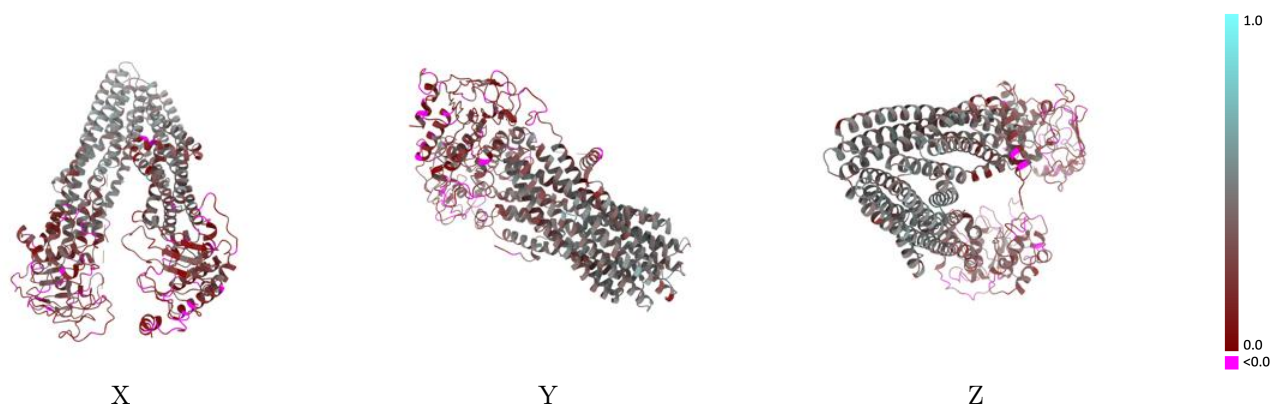
This section contains information regarding the fit between EMDB map EMD-61827 and PDB model 9JUU. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



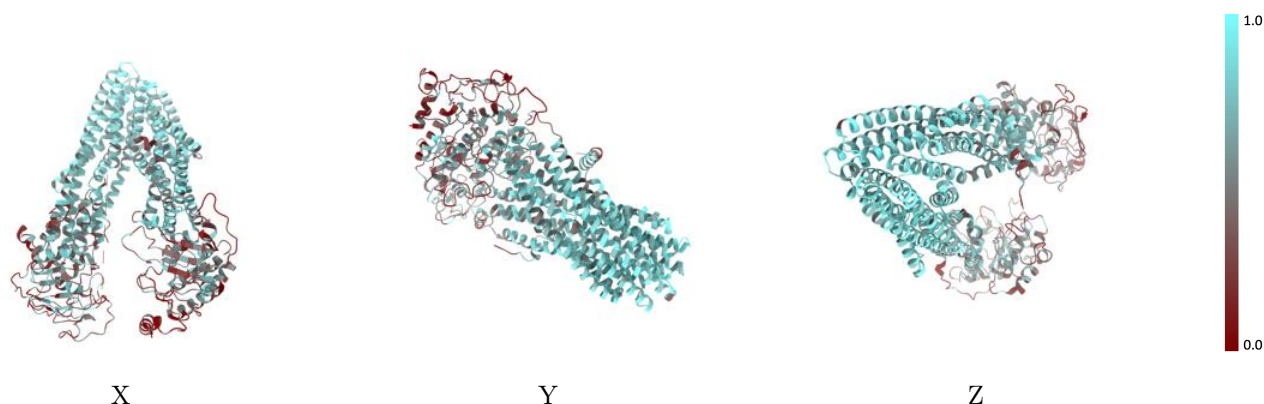
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



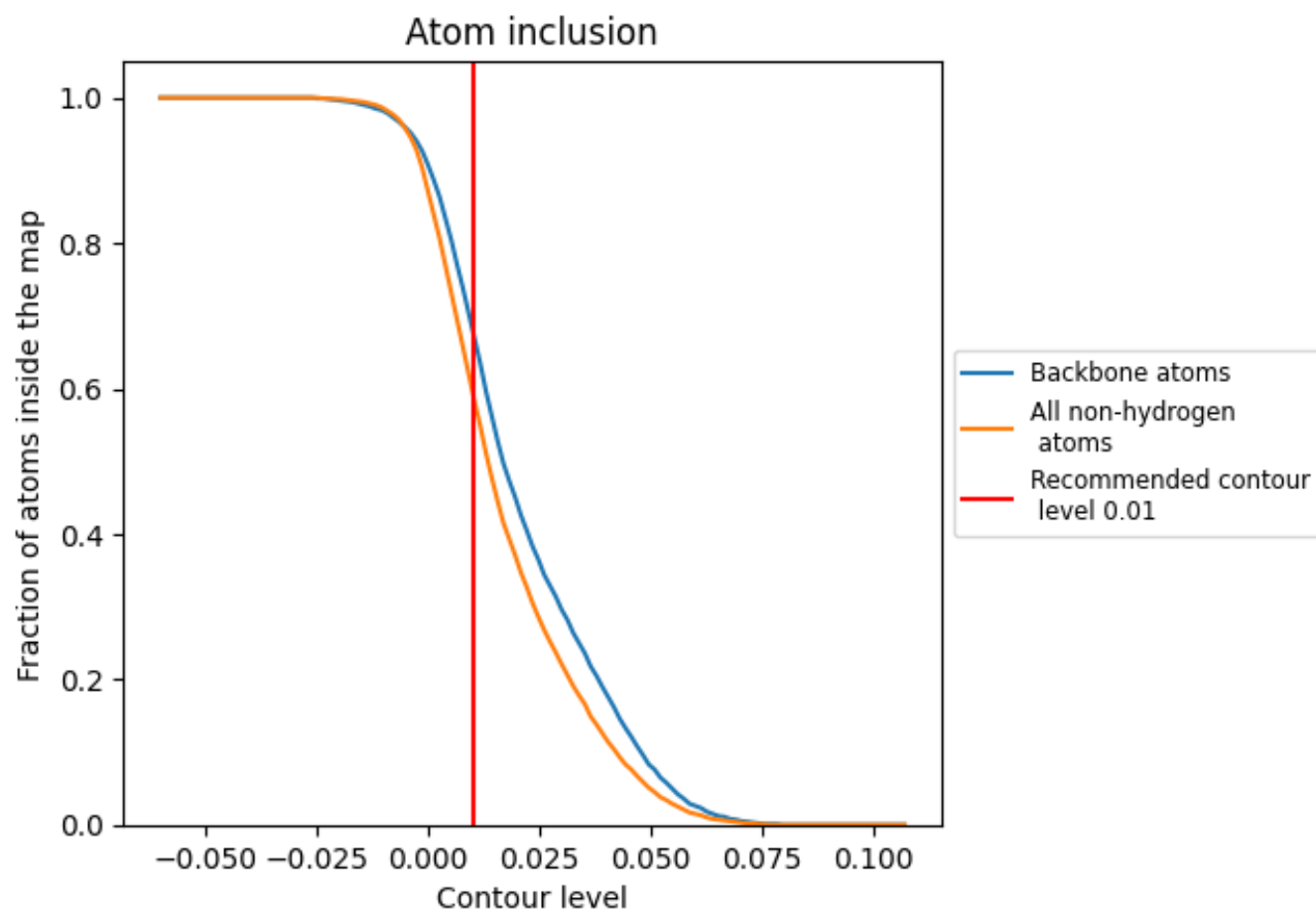
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).

9.4 Atom inclusion [i](#)



At the recommended contour level, 68% of all backbone atoms, 59% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5910	<div></div> 0.2970
A	<div></div> 0.5910	<div></div> 0.2970

