



Full wwPDB EM Validation Report ⓘ

Mar 12, 2025 – 02:22 PM JST

PDB ID : 9JUL
EMDB ID : EMD-61829
Title : Structure of Arabidopsis thaliana ABCB1 with brassinolide bound in the inward-facing conformation
Authors : Chen, Q.; Su, N.; Guo, J.
Deposited on : 2024-10-08
Resolution : 3.10 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
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

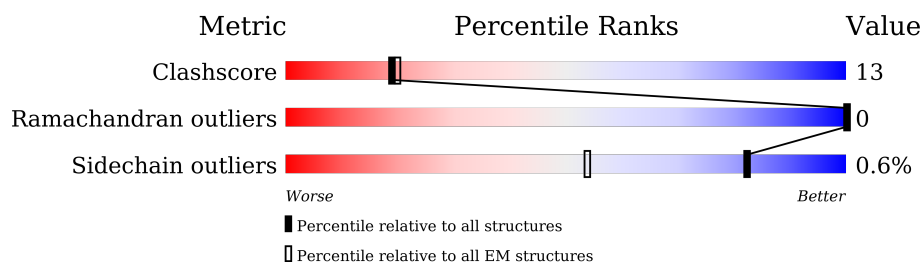
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 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>25%</div> <div>63%</div> <div>24%</div> <div>12%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9005 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:

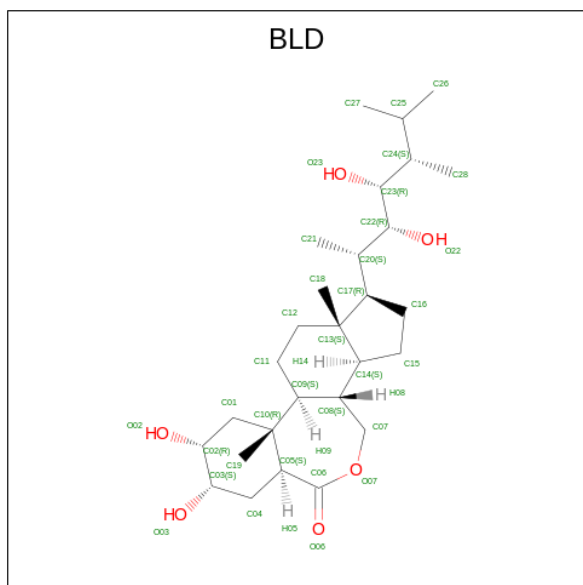
Chain	Residue	Modelled	Actual	Comment	Reference
A	-40	MET	-	initiating methionine	UNP Q9ZR72
A	-39	ASP	-	expression tag	UNP Q9ZR72
A	-38	TYR	-	expression tag	UNP Q9ZR72
A	-37	LYS	-	expression tag	UNP Q9ZR72
A	-36	ASP	-	expression tag	UNP Q9ZR72
A	-35	ASP	-	expression tag	UNP Q9ZR72
A	-34	ASP	-	expression 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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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

- Molecule 2 is Brassinolide (three-letter code: BLD) (formula: $C_{28}H_{48}O_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
2	A	1	Total	C	O	0
			34	28	6	

G989	L1057	G1126	A1192	D1263
R994	V1058	A1127	A1193	G1264
S996	G1059	T1128	D1194	I1265
E998	P1060	E1129	A1195	Y1266
D1001	S1061	A1130	E1196	A1267
R1002	G1062	E1131	S1197	R1268
K1003	C1063	I1132	E1198	M1269
T1004	G1064	T1133	R1199	I1260
E1005	K1065	Q1134	S1200	Q1261
	S1066	A1135	V1201	L1262
	S1067	A1136	Q1202	Q1263
	V1068	T1137	E1203	R1264
	I1069	L1138	A1204	PHE
	S1070	A1139	L1205	THR
	L1071	S1140	D1206	HIS
	Q1073	A1141	Q1207	THR
		H1142	A1208	GLN
	P1078	K1143	C1209	VAL
	S1079	F1144	S1210	ILE
		I1145	G1211	GLY
	R1082	S1146	R1212	MET
	V1083	A1147	T1213	THR
	M1084	L1148		GLY
	I1085	P1149	V1216	SER
	D1086	E1150	V1217	SER
	G1087	G1151	A1218	SER
	K1088	Y1152	H1219	ARG
	I1089		R1220	VAL
	I1090		L1221	LYS
	R1091		S1222	GLU
	K1092		T1223	ASP
	Y1093		I1224	ALA
	N1094		R1225	
	L1095		N1226	
	K1096		A1227	
	A1097		H1228	
	I1098		V1229	
	R1099		I1230	
	K1100		A1231	
			V1232	
	Q1107		I1233	
	E1108		D1234	
	P1109		D1235	
	C1110		G1236	
	L1111		K1237	
	F1112		V1238	
	G1113		A1239	
	T1114		E1240	
	L1115		Q1241	
	I1116		G1242	
	Y1117		S1243	
	E1118		H1244	
	N1119		S1245	
	I1120		H1246	
	A1121		L1247	
	Y1122		L1248	
	G1123		K1249	
	H1124		N1250	
	E1125		H1251	
			P1252	

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	197075	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	0.078	Depositor
Minimum map value	-0.044	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.014	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](#)

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

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	0.34	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	227	0
2	A	34	0	48	2	0
All	All	9005	0	9164	229	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 13.

All (229) 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:216:SER:O	1:A:220:GLN:HG2	1.56	1.04
1:A:199:PRO:O	1:A:203:VAL:HG23	1.71	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1021:ARG:HG3	1:A:1051:ALA:HB1	1.62	0.81
1:A:220:GLN:HE21	1:A:220:GLN:HA	1.50	0.76
1:A:1050:ARG:HB3	1:A:1053:LYS:HD3	1.69	0.72
1:A:1032:SER:HA	1:A:1040:GLN:HA	1.73	0.70
1:A:1130:ALA:O	1:A:1134:GLN:NE2	2.26	0.69
1:A:547:GLU:HB2	1:A:551:ARG:HH21	1.57	0.69
1:A:1123:GLY:O	1:A:1179:ARG:NH1	2.28	0.67
1:A:798:GLU:O	1:A:801:ARG:HG2	1.96	0.65
1:A:466:LEU:HD21	1:A:519:ARG:HH11	1.61	0.65
1:A:428:LEU:HA	1:A:433:ASP:HA	1.78	0.65
1:A:1033:TYR:HD1	1:A:1034:PRO:HD2	1.61	0.64
1:A:1084:MET:SD	1:A:1087:GLY:N	2.70	0.64
1:A:528:LEU:H	1:A:555:GLY:HA2	1.62	0.63
1:A:220:GLN:HA	1:A:220:GLN:NE2	2.14	0.62
1:A:513:GLN:O	1:A:517:ILE:HD12	2.00	0.62
1:A:722:SER:HB3	1:A:964:ARG:HA	1.81	0.62
1:A:63:ARG:NH1	1:A:67:ASP:OD1	2.33	0.61
1:A:1182:GLU:HA	1:A:1212:ARG:HD2	1.81	0.61
1:A:811:ASN:O	1:A:814:ARG:HG2	2.01	0.60
1:A:401:LEU:HD11	1:A:577:LEU:HB2	1.84	0.60
1:A:1021:ARG:HG3	1:A:1051:ALA:CB	2.32	0.60
1:A:370:LEU:HD23	1:A:391:LEU:HB2	1.85	0.59
1:A:464:ILE:HG22	1:A:522:LEU:HD11	1.84	0.59
1:A:782:SER:OG	1:A:786:LYS:NZ	2.34	0.59
1:A:1014:THR:H	1:A:1094:ASN:HD21	1.51	0.59
1:A:1052:GLY:H	1:A:1213:THR:HB	1.66	0.59
1:A:367:LEU:HA	1:A:395:ALA:HB2	1.84	0.59
1:A:1140:SER:O	1:A:1170:ARG:NH1	2.36	0.59
1:A:569:ARG:NH2	1:A:588:HIS:O	2.36	0.58
1:A:1198:GLU:O	1:A:1202:GLN:HG3	2.03	0.58
1:A:123:ARG:NH1	1:A:155:ALA:HA	2.18	0.58
1:A:198:VAL:HG11	1:A:318:GLY:HA3	1.85	0.58
1:A:774:LYS:O	1:A:778:GLU:HG3	2.04	0.58
1:A:368:VAL:HG13	1:A:371:LYS:HZ1	1.69	0.57
1:A:116:GLU:OE1	1:A:159:GLN:NE2	2.37	0.57
1:A:227:GLY:O	1:A:231:GLU:HG3	2.05	0.57
1:A:1187:ASP:OD1	1:A:1188:GLU:N	2.36	0.57
1:A:517:ILE:O	1:A:521:MET:HE2	2.04	0.56
1:A:197:VAL:CG1	1:A:315:MET:HE3	2.35	0.56
1:A:1027:LYS:HA	1:A:1046:SER:HA	1.88	0.56
1:A:1115:THR:N	1:A:1118:GLU:OE2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ASP:O	1:A:143:ARG:NH1	2.39	0.55
1:A:197:VAL:CG1	1:A:315:MET:CE	2.84	0.55
1:A:957:SER:OG	1:A:958:ASP:N	2.39	0.55
1:A:1190:THR:O	1:A:1220:ARG:NH1	2.38	0.55
1:A:83:MET:HG3	1:A:953:LYS:HD2	1.89	0.55
1:A:290:LEU:HD13	1:A:724:VAL:HG21	1.88	0.55
1:A:197:VAL:HG12	1:A:315:MET:CE	2.37	0.55
1:A:122:MET:HE1	1:A:154:ASP:HB2	1.89	0.54
1:A:393:VAL:HB	1:A:556:ARG:HD3	1.89	0.54
1:A:154:ASP:OD1	1:A:155:ALA:N	2.40	0.54
1:A:677:LEU:HD11	1:A:994:ARG:HH12	1.72	0.54
1:A:677:LEU:HD21	1:A:994:ARG:HH22	1.72	0.54
1:A:122:MET:CE	1:A:154:ASP:HB2	2.38	0.54
1:A:58:LEU:HB2	1:A:59:PRO:HD3	1.89	0.54
1:A:517:ILE:HG22	1:A:521:MET:HE2	1.90	0.54
1:A:1145:ILE:HD13	1:A:1152:TYR:HE1	1.71	0.54
1:A:696:GLU:OE1	1:A:771:ASN:ND2	2.34	0.53
1:A:1170:ARG:HH21	1:A:1200:SER:HB2	1.74	0.53
1:A:1225:ARG:NH2	1:A:1245:SER:OG	2.41	0.53
1:A:148:VAL:O	1:A:151:ILE:HG22	2.08	0.53
1:A:273:MET:HA	1:A:759:THR:HG22	1.89	0.53
1:A:1142:HIS:HA	1:A:1145:ILE:HB	1.90	0.53
1:A:245:GLY:HA2	1:A:248:ARG:HD3	1.91	0.53
1:A:245:GLY:HA3	1:A:1122:TYR:OH	2.09	0.53
1:A:385:ILE:HG21	1:A:408:GLY:HA3	1.90	0.52
1:A:1031:PHE:HD1	1:A:1078:PRO:HA	1.73	0.52
1:A:115:GLY:HA3	1:A:163:SER:HB3	1.92	0.52
1:A:714:SER:HB2	1:A:830:LEU:HD13	1.91	0.52
1:A:111:TRP:HB3	1:A:163:SER:O	2.10	0.52
1:A:487:SER:OG	1:A:511:GLN:NE2	2.33	0.52
1:A:715:ALA:HB2	1:A:830:LEU:HD12	1.92	0.52
1:A:116:GLU:HG3	1:A:916:LEU:HD12	1.92	0.52
1:A:386:LEU:HD13	1:A:389:PHE:CZ	2.44	0.52
1:A:795:GLU:HA	1:A:801:ARG:HH22	1.74	0.52
1:A:481:ARG:HH12	1:A:489:ILE:HD11	1.74	0.51
1:A:485:ALA:H	1:A:514:ARG:HH12	1.56	0.51
1:A:773:THR:CG2	1:A:814:ARG:HB3	2.40	0.51
1:A:197:VAL:HG11	1:A:315:MET:HE3	1.93	0.51
1:A:565:LEU:HA	1:A:568:ILE:HD12	1.94	0.51
1:A:816:ALA:HB2	1:A:989:GLY:HA3	1.92	0.50
1:A:1059:GLY:HA3	1:A:1233:ILE:HG22	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1089:ASP:OD1	1:A:1090:ILE:N	2.44	0.50
1:A:850:VAL:HA	1:A:853:PHE:HB2	1.94	0.50
1:A:123:ARG:HH12	1:A:155:ALA:HA	1.77	0.50
1:A:208:HIS:CD2	1:A:274:GLY:HA3	2.47	0.50
1:A:485:ALA:N	1:A:514:ARG:HH12	2.10	0.50
1:A:1045:LEU:HD21	1:A:1238:VAL:HB	1.93	0.50
1:A:114:SER:O	1:A:118:GLN:HG3	2.12	0.50
1:A:517:ILE:HG22	1:A:521:MET:CE	2.42	0.50
1:A:394:PRO:HD2	1:A:399:ILE:HD11	1.94	0.49
1:A:1031:PHE:CD1	1:A:1078:PRO:HA	2.47	0.49
1:A:540:GLU:O	1:A:543:LYS:HG3	2.11	0.49
1:A:574:VAL:HG12	1:A:574:VAL:O	2.13	0.49
1:A:1069:ILE:HB	1:A:1185:LEU:HD13	1.93	0.49
1:A:478:GLU:O	1:A:482:VAL:HG12	2.11	0.49
1:A:119:THR:HG21	1:A:159:GLN:HB2	1.95	0.49
1:A:488:PHE:O	1:A:492:LEU:HB2	2.13	0.49
1:A:500:VAL:HG23	1:A:501:GLY:H	1.78	0.49
1:A:995:SER:O	1:A:998:GLU:HG3	2.13	0.48
1:A:460:ILE:HA	1:A:463:ASN:HB2	1.96	0.48
1:A:1070:SER:HA	1:A:1073:GLN:HE21	1.77	0.48
1:A:1186:LEU:HB2	1:A:1216:VAL:HG22	1.94	0.48
1:A:402:VAL:HG13	1:A:561:ILE:HA	1.95	0.48
1:A:468:ARG:HE	1:A:471:ALA:HB2	1.79	0.48
1:A:544:LEU:O	1:A:547:GLU:HG2	2.13	0.48
1:A:1010:ASP:OD1	1:A:1010:ASP:N	2.46	0.48
1:A:1026:LEU:HD13	1:A:1047:LEU:HD23	1.95	0.48
1:A:372:ASN:N	1:A:390:CYS:SG	2.77	0.48
1:A:75:ASN:HB2	1:A:82:MET:HB2	1.95	0.48
1:A:1031:PHE:CZ	1:A:1034:PRO:HD3	2.48	0.48
1:A:439:LEU:HA	1:A:442:LEU:HB3	1.96	0.48
1:A:57:SER:HB2	1:A:100:ILE:HD11	1.96	0.48
1:A:1057:LEU:HB2	1:A:1217:VAL:HG12	1.96	0.47
1:A:193:VAL:HG12	1:A:289:LEU:CD2	2.44	0.47
1:A:453:PRO:HG2	1:A:512:LYS:HG3	1.95	0.47
1:A:258:LYS:O	1:A:262:LYS:HG2	2.15	0.47
1:A:85:GLU:O	1:A:88:LYS:HG2	2.13	0.47
1:A:275:LEU:HD12	1:A:762:HIS:NE2	2.30	0.47
1:A:397:LYS:HD2	1:A:397:LYS:HA	1.78	0.47
1:A:527:ILE:HD12	1:A:556:ARG:HH11	1.80	0.47
1:A:368:VAL:HG23	1:A:393:VAL:H	1.79	0.47
1:A:677:LEU:HD11	1:A:994:ARG:NH1	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LEU:HD13	1:A:945:LEU:HD23	1.97	0.47
1:A:1045:LEU:HD22	1:A:1047:LEU:HB2	1.96	0.47
1:A:234:VAL:HG11	1:A:799:SER:H	1.80	0.47
1:A:459:SER:HA	1:A:499:GLN:HA	1.97	0.47
1:A:460:ILE:O	1:A:464:ILE:HG12	2.15	0.47
1:A:485:ALA:O	1:A:514:ARG:NH2	2.45	0.47
1:A:165:LYS:HA	1:A:165:LYS:HD3	1.71	0.46
1:A:1114:THR:HA	1:A:1155:TYR:HD1	1.80	0.46
1:A:554:ILE:O	1:A:556:ARG:HG3	2.16	0.46
1:A:1169:GLN:NE2	1:A:1189:ALA:O	2.49	0.46
1:A:709:ILE:HD12	1:A:757:PHE:CE2	2.51	0.46
1:A:1021:ARG:HH21	1:A:1213:THR:HG22	1.80	0.46
1:A:290:LEU:HD11	1:A:721:LEU:HD23	1.97	0.46
1:A:705:VAL:O	1:A:708:VAL:HG22	2.15	0.46
1:A:358:GLY:HA2	1:A:437:LEU:HA	1.98	0.46
1:A:486:HIS:HB3	1:A:490:ILE:HD11	1.98	0.46
1:A:101:TRP:CD1	1:A:931:GLY:HA3	2.51	0.45
1:A:1018:ASP:N	1:A:1018:ASP:OD1	2.47	0.45
1:A:1041:ILE:HD13	1:A:1067:SER:HB2	1.99	0.45
1:A:1174:ALA:O	1:A:1178:VAL:HG23	2.16	0.45
1:A:241:MET:HE1	1:A:789:MET:HG2	1.98	0.45
1:A:1016:VAL:HG22	1:A:1093:TYR:CZ	2.51	0.45
1:A:1042:PHE:CE2	1:A:1045:LEU:HB3	2.52	0.45
1:A:798:GLU:H	1:A:801:ARG:HE	1.65	0.45
1:A:487:SER:OG	1:A:488:PHE:N	2.49	0.44
2:A:1301:BLD:H05	2:A:1301:BLD:H107	1.65	0.44
1:A:290:LEU:HD22	1:A:724:VAL:HG11	1.99	0.44
1:A:727:VAL:HG12	1:A:736:MET:HG3	2.00	0.44
1:A:1069:ILE:O	1:A:1073:GLN:HG2	2.18	0.44
1:A:446:ILE:HG23	1:A:527:ILE:HG23	1.99	0.44
1:A:1049:ALA:HB2	1:A:1055:LEU:HD22	2.00	0.44
1:A:387:ASN:OD1	1:A:388:ASN:N	2.47	0.43
1:A:432:GLN:O	1:A:436:THR:OG1	2.36	0.43
1:A:569:ARG:HH21	1:A:588:HIS:HB2	1.83	0.43
1:A:1028:HIS:H	1:A:1046:SER:HA	1.83	0.43
1:A:197:VAL:HG11	1:A:315:MET:CE	2.47	0.43
1:A:257:LEU:HD12	1:A:778:GLU:HG2	1.99	0.43
1:A:1089:ASP:OD1	1:A:1091:ARG:N	2.31	0.43
1:A:197:VAL:CG1	1:A:315:MET:HE1	2.48	0.43
1:A:403:GLY:H	1:A:409:LYS:HD3	1.83	0.43
1:A:395:ALA:HA	1:A:556:ARG:HH21	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1095:LEU:O	1:A:1099:ARG:HG2	2.19	0.43
1:A:492:LEU:HD12	1:A:493:PRO:HD2	2.01	0.43
1:A:963:ILE:HD12	1:A:963:ILE:HA	1.90	0.43
1:A:157:MET:HE3	1:A:337:ALA:HB2	2.00	0.43
1:A:187:VAL:HG12	1:A:190:LEU:H	1.83	0.43
1:A:409:LYS:HE3	1:A:561:ILE:HD12	2.01	0.43
1:A:755:LEU:O	1:A:759:THR:HG23	2.18	0.42
1:A:1144:PHE:CD2	1:A:1167:GLN:HG2	2.54	0.42
1:A:27:VAL:H	1:A:339:ALA:HB2	1.83	0.42
1:A:1240:GLU:HB3	1:A:1247:LEU:HD13	2.01	0.42
1:A:397:LYS:O	1:A:556:ARG:HG2	2.20	0.42
1:A:1193:LEU:HD23	1:A:1197:SER:HB3	2.01	0.42
1:A:566:SER:O	1:A:570:LYS:HG2	2.19	0.42
1:A:263:LEU:HA	1:A:266:LYS:HD3	2.02	0.42
1:A:475:GLU:O	1:A:478:GLU:HG3	2.20	0.42
1:A:802:ILE:O	1:A:805:ARG:HG2	2.20	0.42
1:A:149:PHE:O	1:A:153:THR:HG23	2.20	0.42
1:A:197:VAL:HG12	1:A:315:MET:HE1	2.01	0.42
1:A:241:MET:HE2	1:A:241:MET:HB2	1.85	0.42
1:A:568:ILE:HG23	1:A:574:VAL:HG11	2.02	0.42
1:A:1232:VAL:HG21	1:A:1255:ILE:HG21	2.01	0.42
1:A:237:ILE:HG13	1:A:241:MET:HE2	2.02	0.42
1:A:433:ASP:CG	1:A:434:LEU:H	2.23	0.42
1:A:489:ILE:HG13	1:A:490:ILE:N	2.34	0.42
1:A:1026:LEU:HB2	1:A:1047:LEU:HB3	2.01	0.42
1:A:240:VAL:HG12	1:A:246:GLU:OE1	2.20	0.41
1:A:1033:TYR:HB3	1:A:1036:ARG:O	2.20	0.41
1:A:35:PHE:HE1	1:A:125:LYS:HD3	1.85	0.41
1:A:492:LEU:HD21	1:A:498:THR:HG21	2.01	0.41
1:A:518:ALA:HA	1:A:521:MET:HE3	2.01	0.41
1:A:716:PHE:CE2	1:A:750:LEU:HD12	2.55	0.41
1:A:811:ASN:OD1	1:A:812:ASN:N	2.51	0.41
1:A:453:PRO:HD3	1:A:513:GLN:NE2	2.35	0.41
1:A:1243:SER:HB3	1:A:1246:HIS:HB3	2.01	0.41
1:A:35:PHE:CD2	1:A:122:MET:HB2	2.56	0.41
1:A:246:GLU:HG3	1:A:785:LEU:HB3	2.02	0.41
1:A:1056:ALA:HB1	1:A:1224:ILE:HG22	2.01	0.41
2:A:1301:BLD:H121	2:A:1301:BLD:H23	1.77	0.41
1:A:367:LEU:HD13	1:A:395:ALA:HB2	2.02	0.41
1:A:474:VAL:O	1:A:477:GLU:HG3	2.21	0.41
1:A:489:ILE:HG13	1:A:490:ILE:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:560:ILE:HD11	1:A:564:ARG:H	1.84	0.41
1:A:373:VAL:HG11	1:A:386:LEU:HD12	2.02	0.41
1:A:1068:VAL:O	1:A:1072:ILE:HG12	2.20	0.41
1:A:237:ILE:HD12	1:A:237:ILE:HA	1.94	0.41
1:A:572:ASP:OD1	1:A:573:LEU:N	2.53	0.41
1:A:1232:VAL:O	1:A:1238:VAL:HA	2.21	0.41
1:A:218:LYS:O	1:A:221:GLU:HG3	2.20	0.41
1:A:369:GLU:HB2	1:A:428:LEU:O	2.21	0.41
1:A:1116:ILE:O	1:A:1120:ILE:HG12	2.21	0.41
1:A:1129:GLU:O	1:A:1133:ILE:HG12	2.20	0.41
1:A:193:VAL:HG12	1:A:289:LEU:HD21	2.03	0.40
1:A:374:ASP:N	1:A:374:ASP:OD1	2.54	0.40
1:A:930:TYR:O	1:A:934:GLN:HG2	2.22	0.40
1:A:948:ALA:HB1	1:A:962:THR:HB	2.03	0.40
1:A:560:ILE:HG12	1:A:563:HIS:H	1.86	0.40
1:A:348:LYS:HD3	1:A:348:LYS:HA	1.89	0.40
1:A:880:LYS:O	1:A:883:GLN:HG3	2.21	0.40
1:A:197:VAL:HG12	1:A:315:MET:HE3	2.02	0.40
1:A:353:ARG:HA	1:A:439:LEU:HD21	2.03	0.40
1:A:418:ARG:HH12	1:A:435:LYS:HA	1.86	0.40
1:A:909:THR:O	1:A:913:GLU:HG2	2.21	0.40
1:A:1052:GLY:N	1:A:1213:THR:HB	2.35	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%)	1100 (95%)	61 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	949/1087 (87%)	943 (99%)	6 (1%)	84	91

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	543	LYS
1	A	1036	ARG
1	A	1042	PHE
1	A	1043	ARG
1	A	1048	ARG
1	A	1069	ILE

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	220	GLN
1	A	236	GLN
1	A	484	ASN
1	A	579	GLN
1	A	1073	GLN
1	A	1134	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [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	$\# Z > 2$
2	BLD	A	1301	-	36,37,37	0.34	0	46,59,59	1.16	2 (4%)

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
2	BLD	A	1301	-	-	8/20/85/85	0/4/4/4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1301	BLD	O07-C06-O06	4.13	123.16	116.72
2	A	1301	BLD	C01-C10-C05	3.63	112.86	107.06

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1301	BLD	C22-C23-C24-C28
2	A	1301	BLD	O22-C22-C23-O23
2	A	1301	BLD	O23-C23-C24-C28
2	A	1301	BLD	O23-C23-C24-C25

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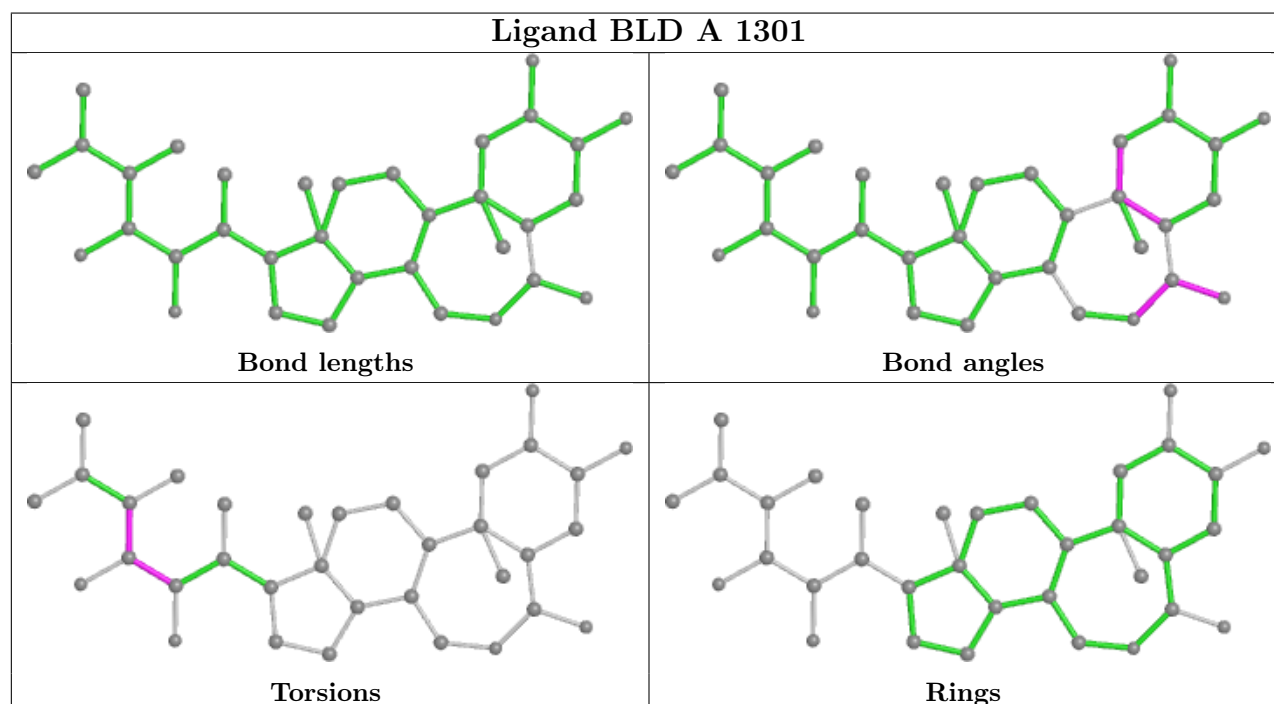
Mol	Chain	Res	Type	Atoms
2	A	1301	BLD	C20-C22-C23-C24
2	A	1301	BLD	C22-C23-C24-C25
2	A	1301	BLD	C20-C22-C23-O23
2	A	1301	BLD	O22-C22-C23-C24

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1301	BLD	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

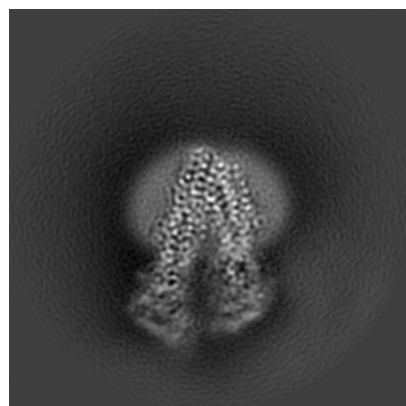
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61829. These allow visual inspection of the internal detail of the map and identification of artifacts.

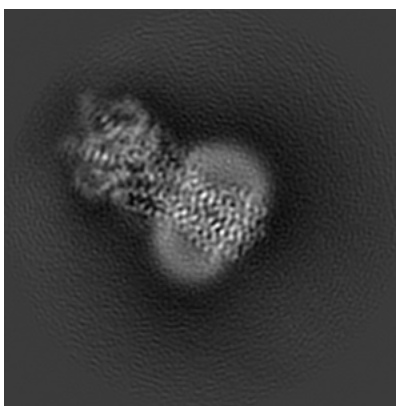
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

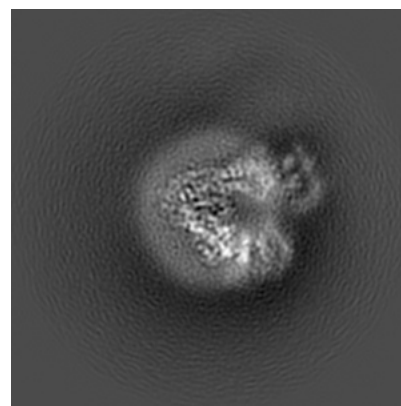
6.1.1 Primary map



X

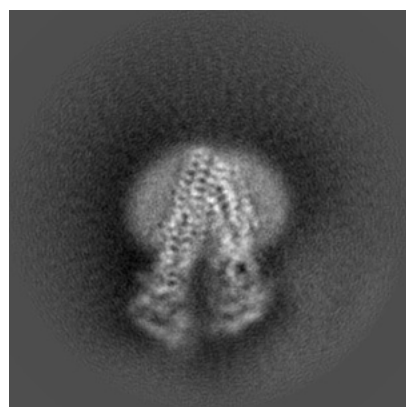


Y

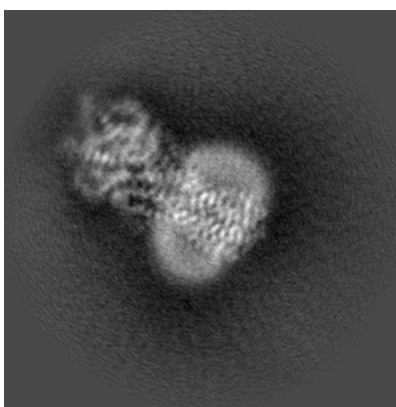


Z

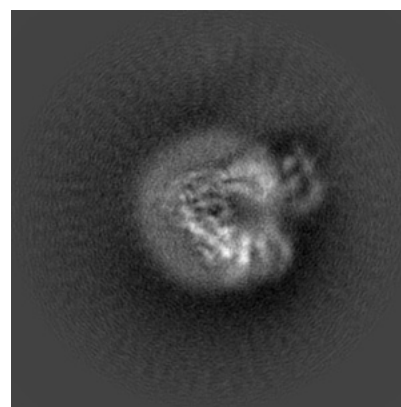
6.1.2 Raw map



X



Y

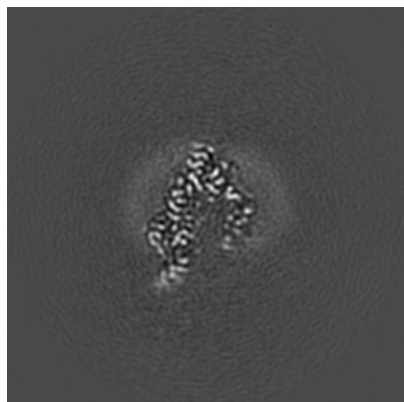


Z

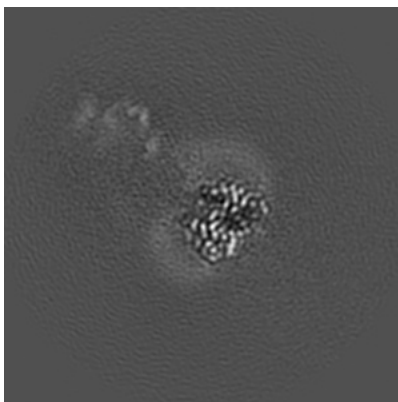
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

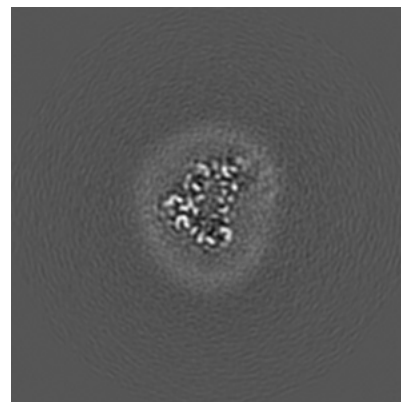
6.2.1 Primary map



X Index: 120

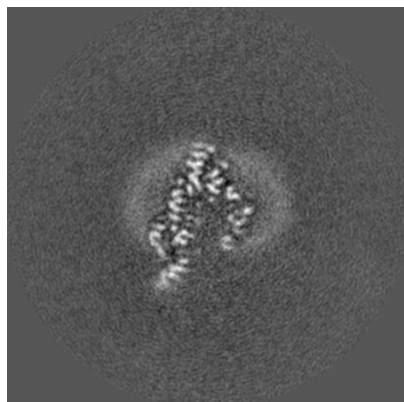


Y Index: 120

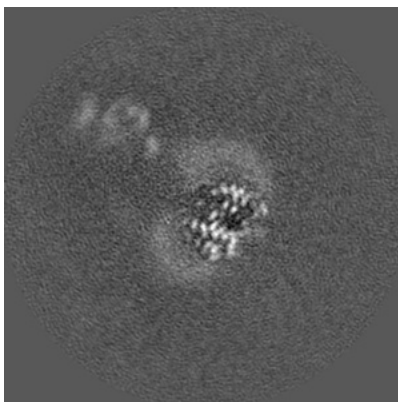


Z Index: 120

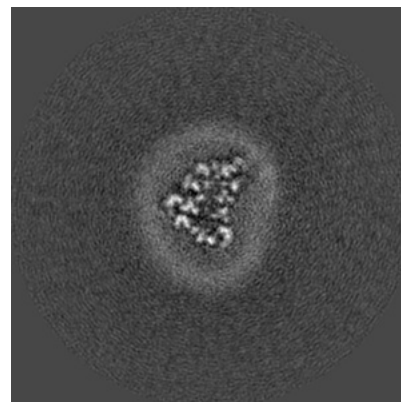
6.2.2 Raw map



X Index: 120



Y Index: 120

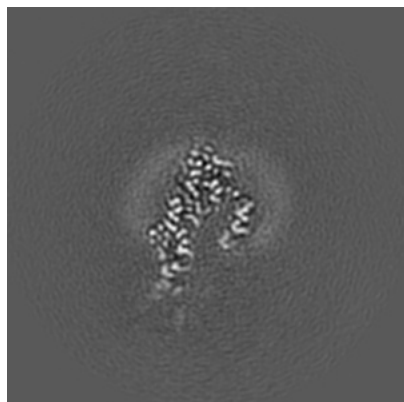


Z Index: 120

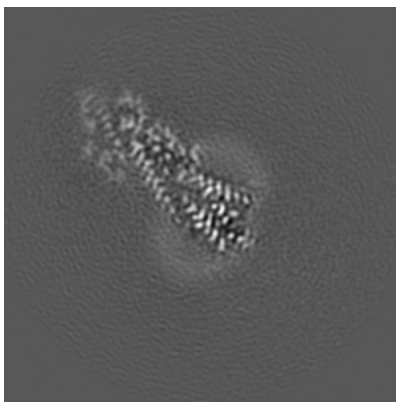
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

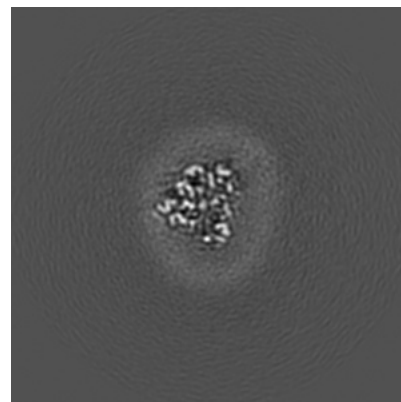
6.3.1 Primary map



X Index: 122

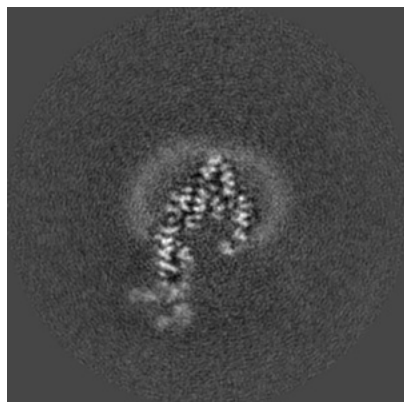


Y Index: 134

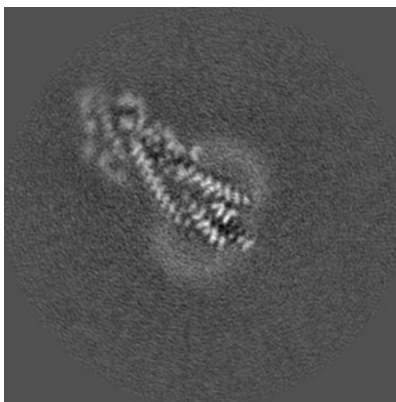


Z Index: 125

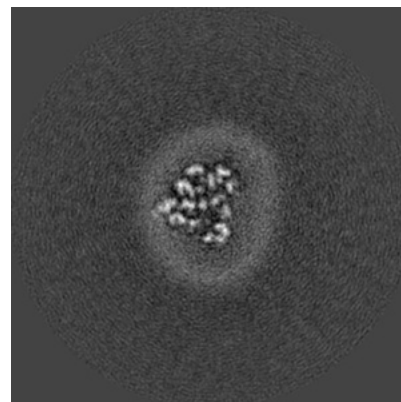
6.3.2 Raw map



X Index: 127



Y Index: 133

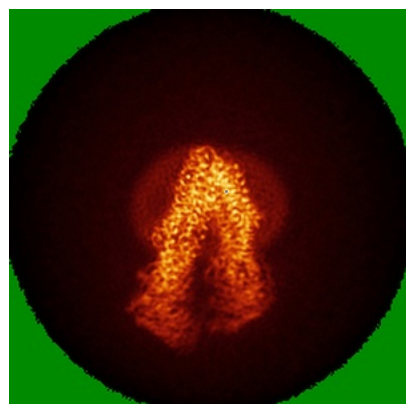


Z Index: 125

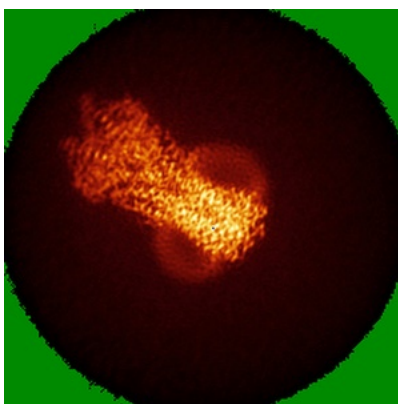
The images above show the largest variance slices of the map in three orthogonal directions.

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

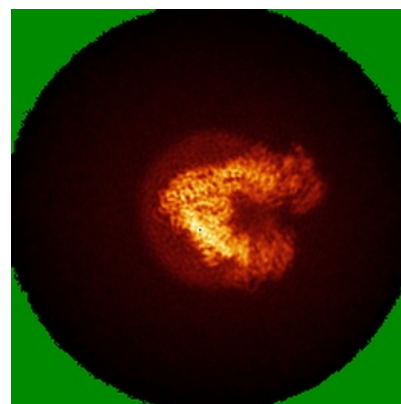
6.4.1 Primary map



X

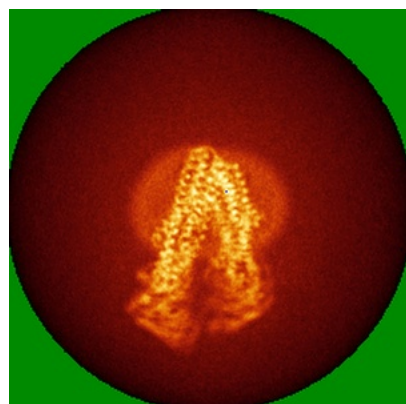


Y

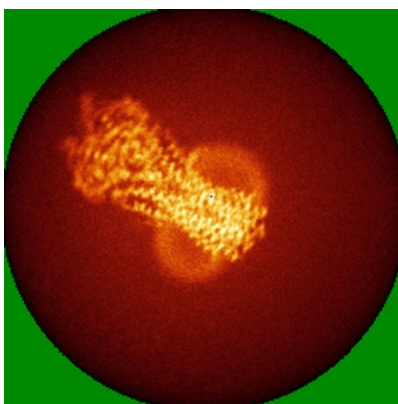


Z

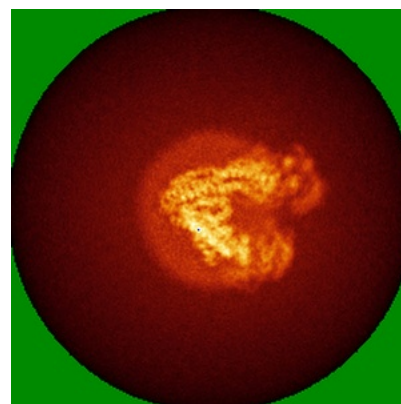
6.4.2 Raw map



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.

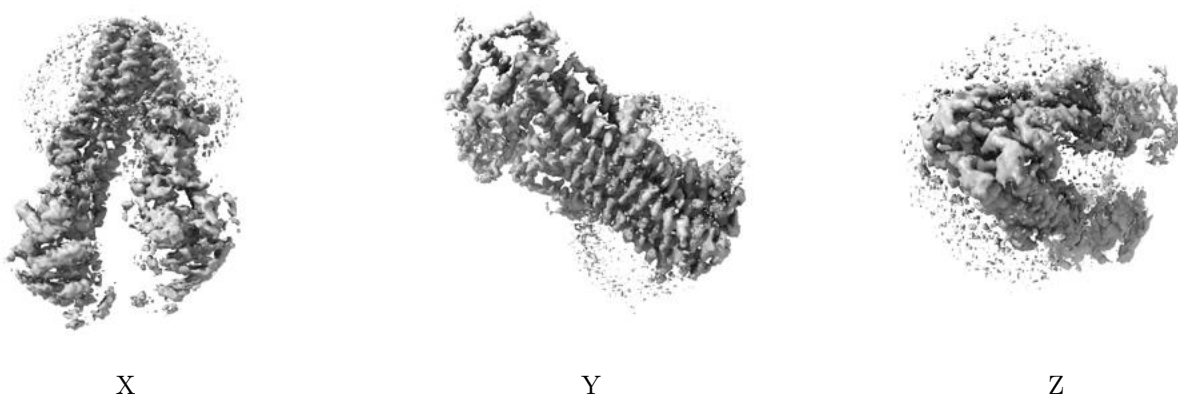
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.014. 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.

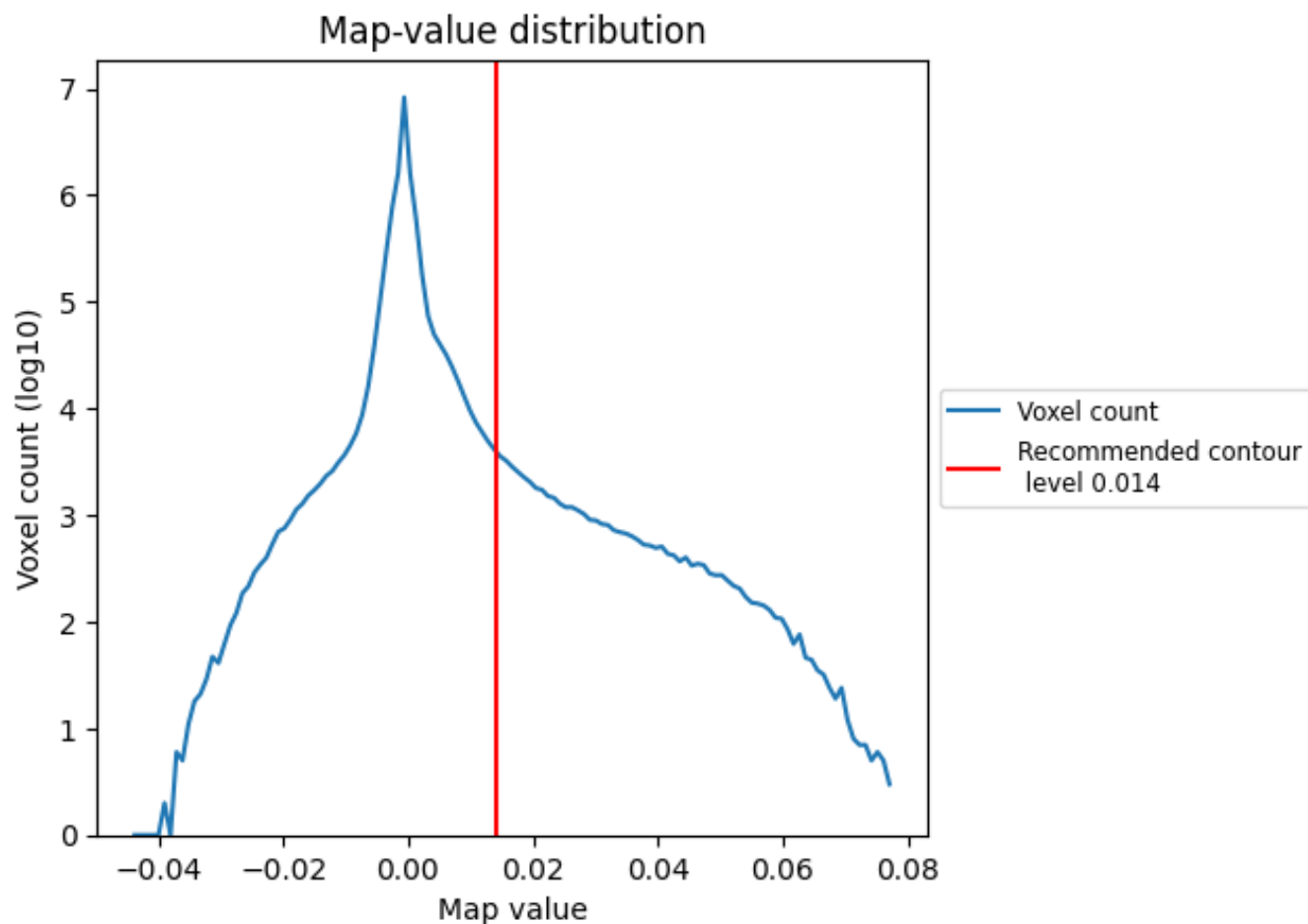
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

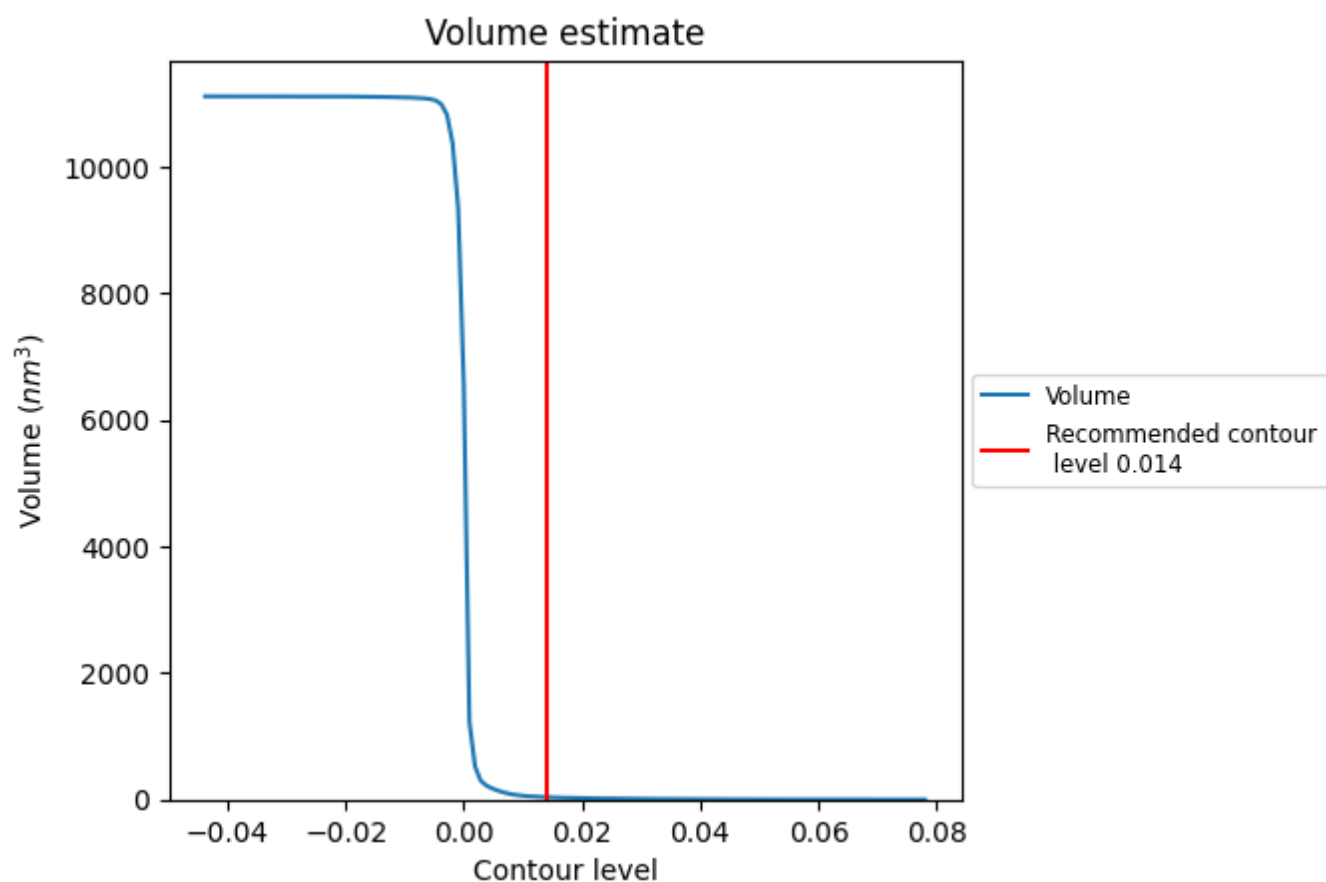
This section contains the results of statistical analysis of the map.

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.

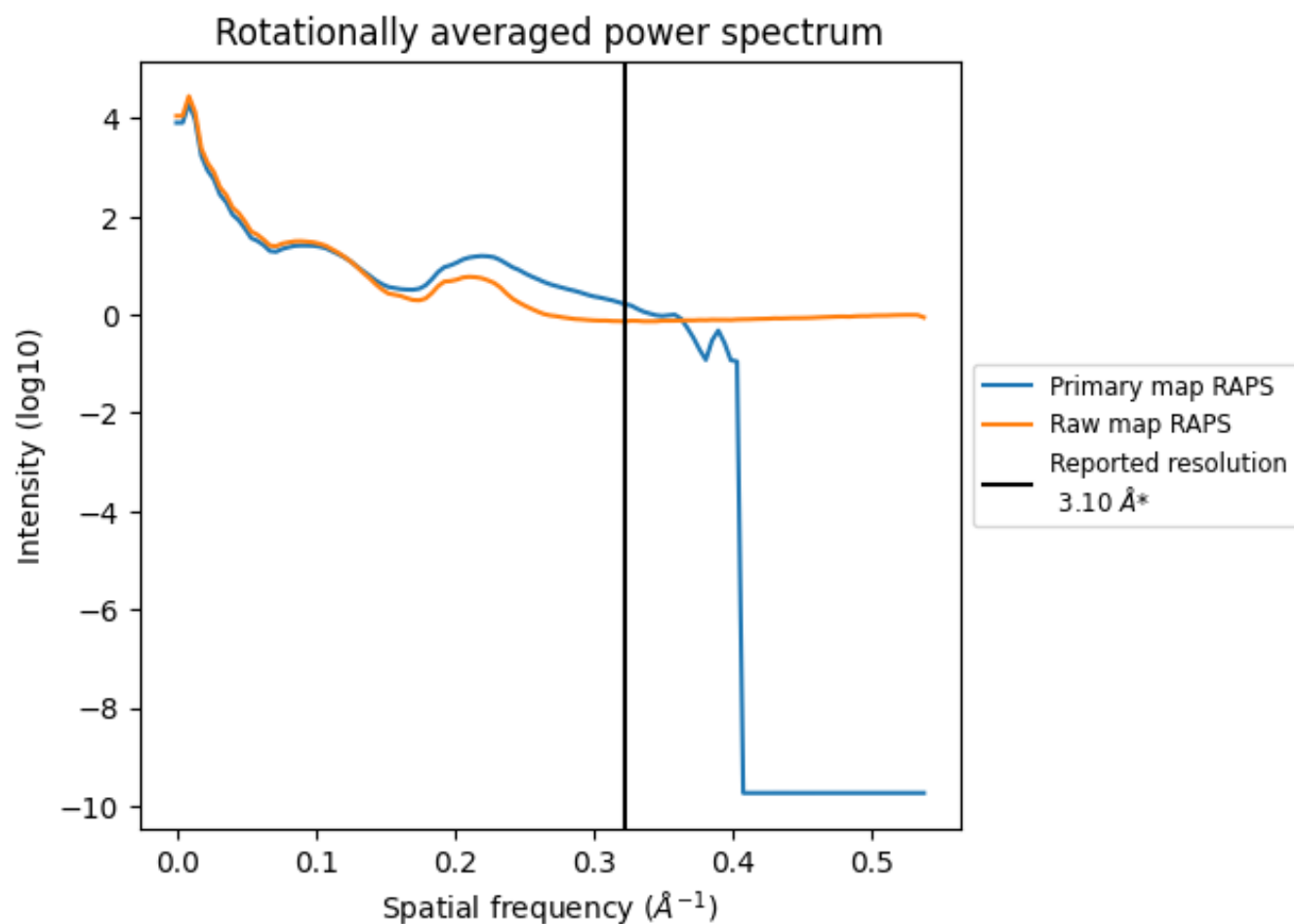
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 37 nm³; this corresponds to an approximate mass of 34 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 ⓘ

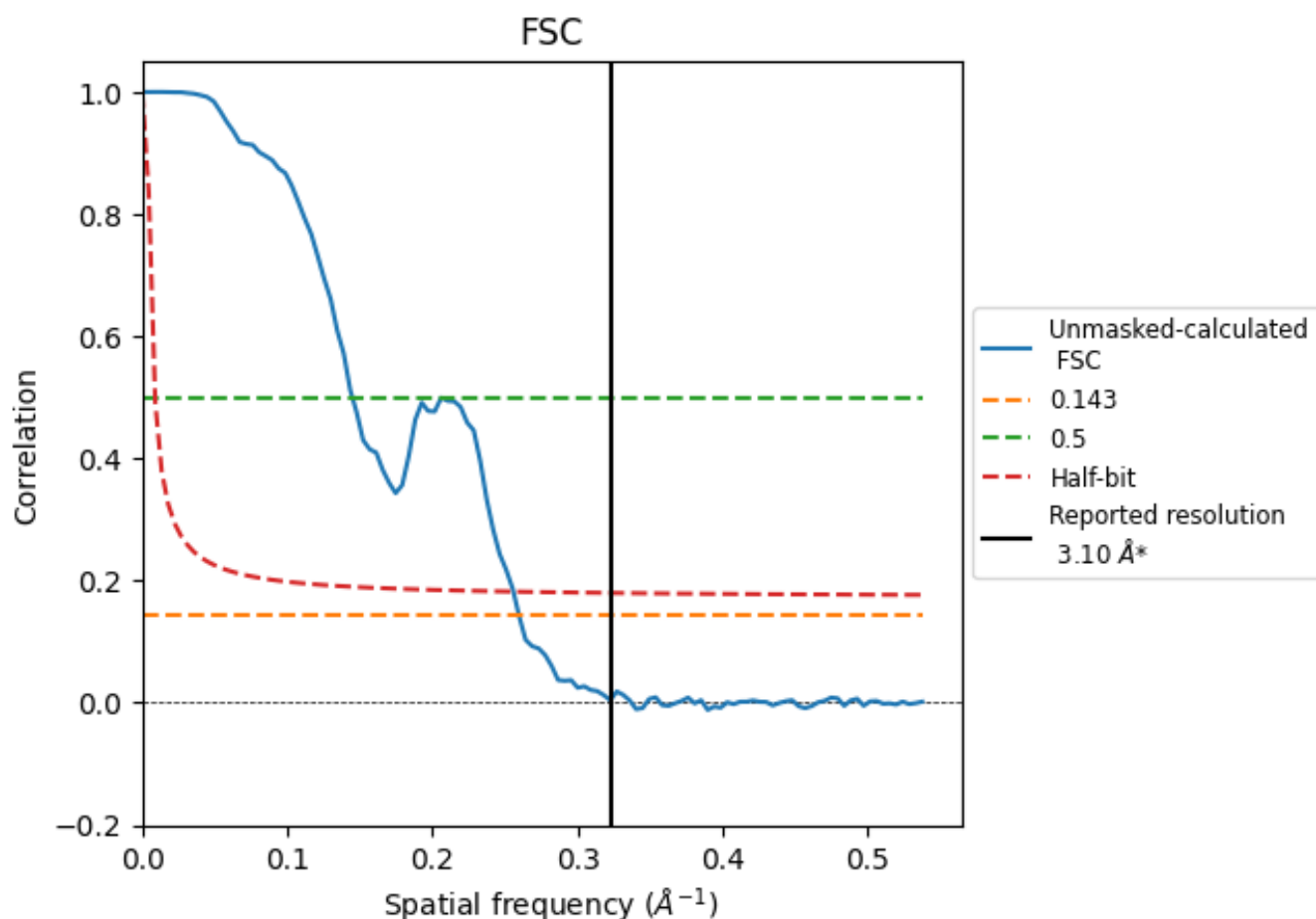


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 0.323 \AA^{-1}

8.2 Resolution estimates [i](#)

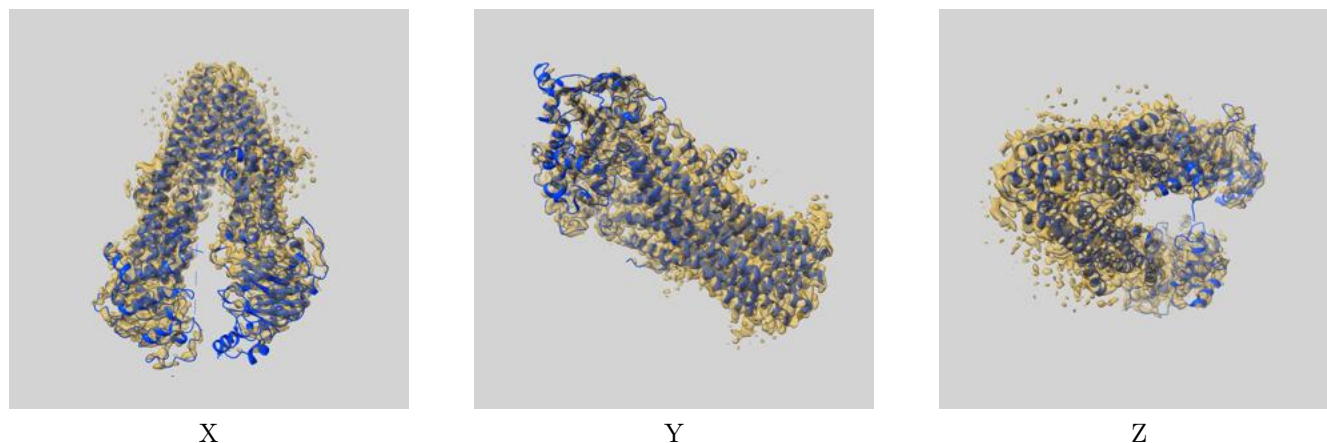
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.85	6.91	3.91

*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 3.85 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

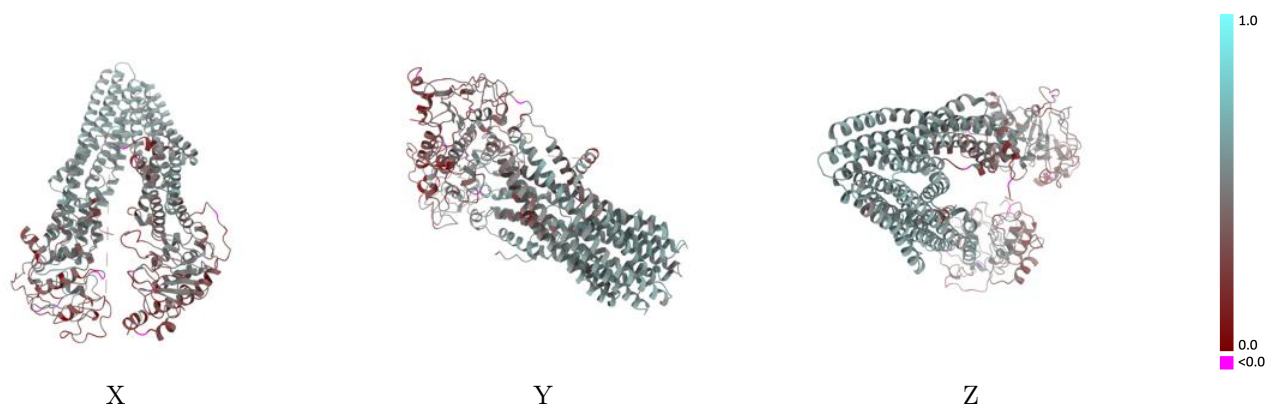
This section contains information regarding the fit between EMDB map EMD-61829 and PDB model 9JUL. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



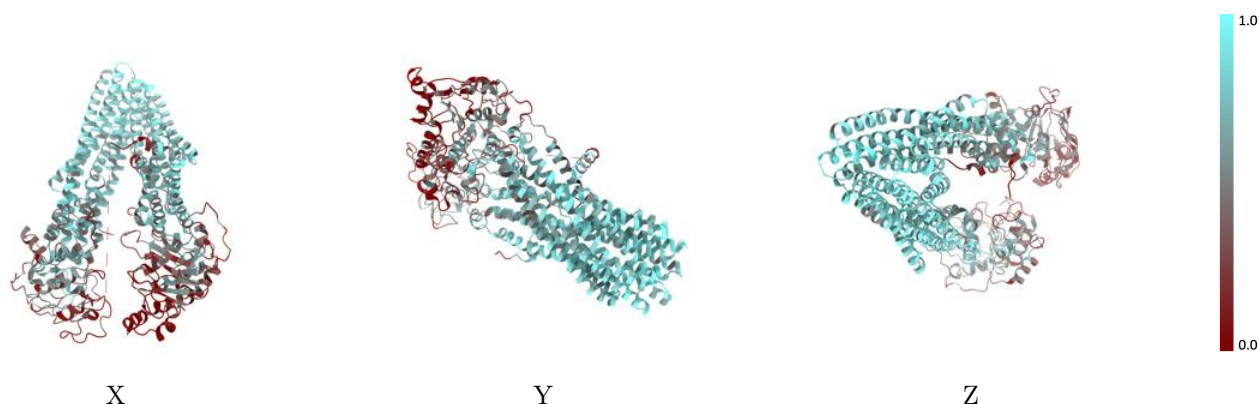
The images above show the 3D surface view of the map at the recommended contour level 0.014 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



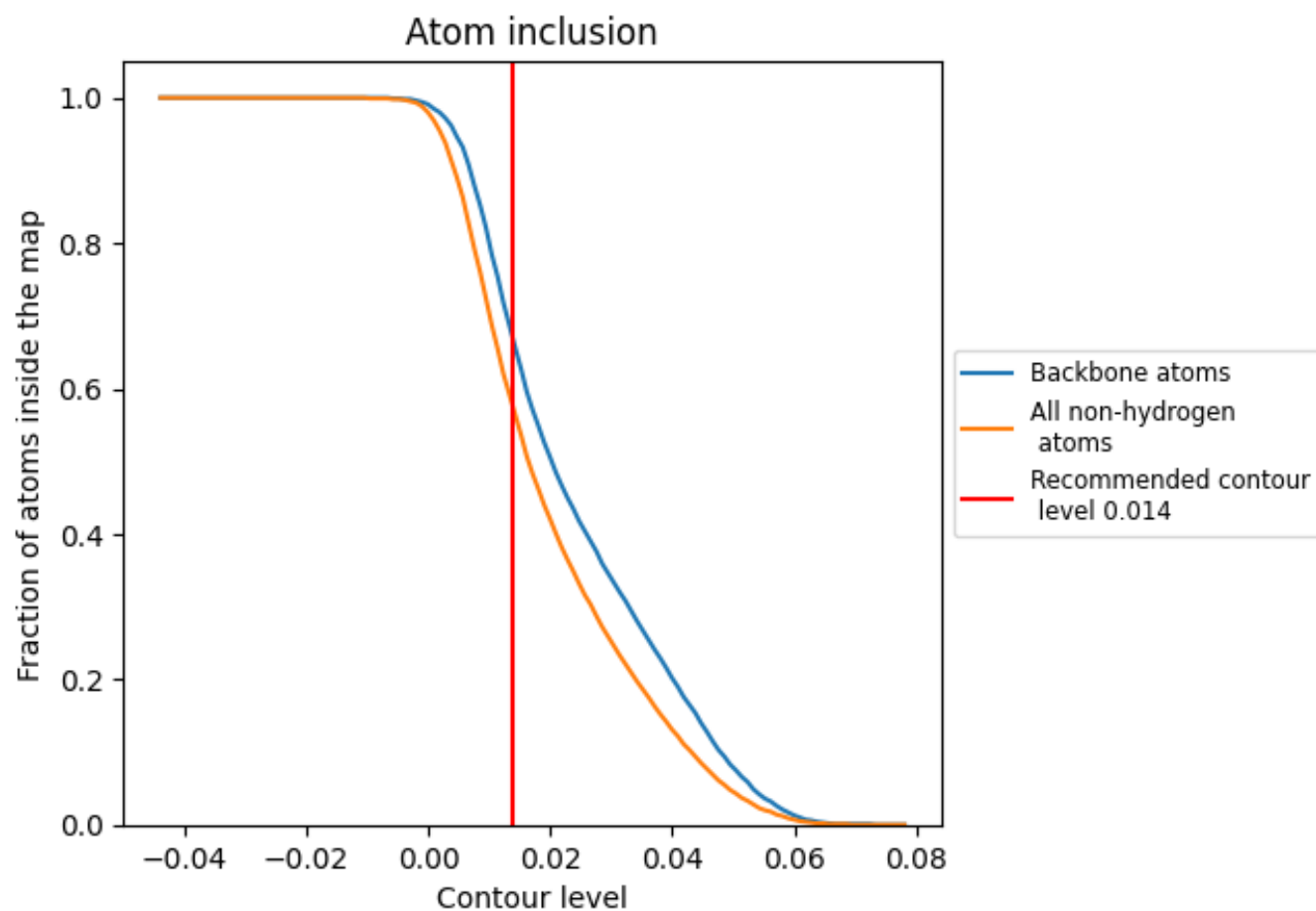
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.014).

9.4 Atom inclusion [i](#)



At the recommended contour level, 66% of all backbone atoms, 57% 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 (0.014) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5720	<div></div> 0.4510
A	<div></div> 0.5720	<div></div> 0.4510

