



Full wwPDB EM Validation Report ⓘ

Mar 25, 2026 – 12:44 AM UTC

PDB ID : 8JQG / pdb_00008jqg
EMDB ID : EMD-36571
Title : Cryo EM map of full length PLC gamma 2
Authors : Shin, Y.-C.; Liao, M.
Deposited on : 2023-06-14
Resolution : 3.72 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

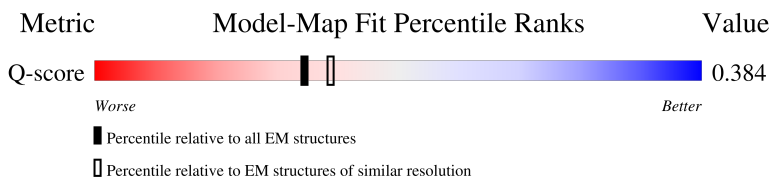
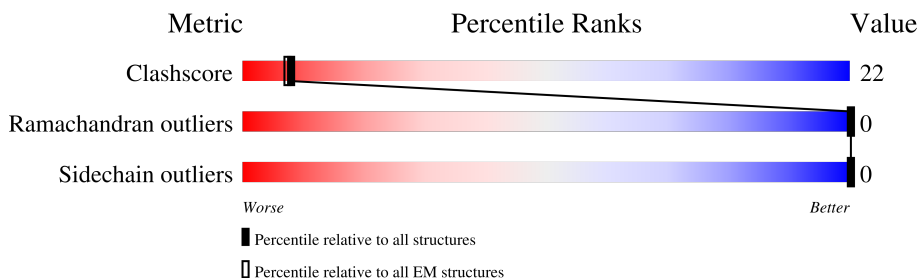
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.72 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10474 (3.22 - 4.22)

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	1265	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8708 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 1-phosphatidylinositol 4,5-bisphosphate phosphodiesterase gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1061	8708	5548	1494	1622	44	0	0

LEU	Y1137	M1043	P939	N868
TYR	E1138	R1044	T940	Q869
ASP			S941	K870
THR	M1146		K942	S871
HIS	F1147	D1049	T943	V872
GLN	L1148	P1050	K944	V873
ASN	A1149	M1051	D945	F874
LEU	H1150		N946	L875
ARG	A1151	I1059	L947	L876
ASN	T1152	L1060		E877
ALA		M1061	P950	P878
ASN	I1155	T1062	D951	K879
ARG		L1063	F952	Q880
ASP	F1162	T1064	R953	
ALA	R1163	V1065	E954	D883
LEU	S1164	K1066	I955	E887
VAL		V1067	R956	A889
LYS	L1167	L1068	S957	T890
GLU		G1069	F958	A889
PHE	Y1171	A1070		D891
SER	S1172	R1071	D964	R892
VAL	E1173			V893
ASN		L1076	I967	E894
GLU	A1178		R968	E895
ASN	S1179	I1080		L896
GLN	L1180	A1081	L975	F897
LEU	L1181	C1082	K976	E898
GLN	V1182	P1083	Y977	V899
LEU		F1084		F900
TYR	E1185	V1085	K980	Q901
GLN	M1186	E1086		S902
GLU	R1187	V1087	V985	I903
LYS	F1188	E1088	F986	R904
CYS	V1189	I1089	P987	E905
ASN	L1190	C1090		I906
LYS	GLU	G1091	D993	T907
SER				V908
ARG	GLU	D1106	N996	LYS
LEU	GLU	N1107		ILE
ARG	GLU	G1108	P999	ASP
GLU	LEU		F1000	THR
LYS	LYS	P1111	R1001	LYS
ARG	TYR	I1112	L1002	GLU
VAL	SER			ASN
SER	SER	Q1117	C1005	ASN
ASN	CYS			MET
SER	ARG	F1122	F1014	LYS
LYS	GLN	E1123		TYR
PHE	LEU	I1124	A1017	TRP
TYR	ARG	Y1125	D1018	GLU
SER	ARG	D1126	K1019	
	ARG	P1127	Y1020	K922
	GLN	N1128	M1021	
	GLU		Q1022	I926
	GLU	F1131	M1023	A927
	LEU	I1132		I928
	ASN	R1133	L1027	L933
	ASN	F1134		
	GLN	V1135	Q1039	
	LEU		P1040	K938
	PHE			

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	189652	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	53.0932	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.982	Depositor
Minimum map value	-1.785	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.075	Depositor
Recommended contour level	0.3	Depositor
Map size ($\text{\AA}$)	237.6, 237.6, 237.6	wwPDB
Map dimensions	216, 216, 216	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	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	0.23	0/8910	0.49	7/12025 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	PHE	N-CA-C	-6.09	105.56	112.87
1	A	218	LEU	N-CA-C	-5.92	105.16	112.38
1	A	540	ARG	N-CA-C	-5.70	105.07	111.28
1	A	577	PHE	N-CA-C	5.58	117.50	108.41
1	A	182	PHE	N-CA-C	-5.22	105.76	111.82
1	A	218	LEU	CA-C-N	-5.06	113.86	120.44
1	A	218	LEU	C-N-CA	-5.06	113.86	120.44

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	8708	0	8605	383	0
All	All	8708	0	8605	383	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 22.

All (383) 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:578:TRP:HE1	1:A:581:GLY:H	1.13	0.92
1:A:491:TRP:HZ2	1:A:868:ASN:HB2	1.40	0.85
1:A:491:TRP:CZ2	1:A:868:ASN:HB2	2.10	0.85
1:A:623:ARG:HG2	1:A:628:GLU:HG2	1.62	0.81
1:A:939:PRO:HA	1:A:955:ILE:HG22	1.63	0.81
1:A:221:PHE:HE1	1:A:249:PHE:CE1	1.97	0.81
1:A:106:LEU:HD11	1:A:347:ARG:HE	1.46	0.81
1:A:22:LEU:HD21	1:A:122:TRP:HA	1.64	0.79
1:A:530:GLU:HG2	1:A:532:TRP:CE2	2.17	0.79
1:A:153:VAL:HG21	1:A:165:GLU:HB3	1.63	0.78
1:A:621:HIS:NE2	1:A:630:ARG:HG2	2.01	0.76
1:A:216:SER:HB3	1:A:1112:ILE:HD11	1.68	0.76
1:A:587:ARG:HE	1:A:589:ARG:HD3	1.52	0.74
1:A:621:HIS:HD2	1:A:630:ARG:N	1.86	0.74
1:A:482:TYR:HB2	1:A:889:ALA:HB3	1.70	0.73
1:A:899:TRP:O	1:A:903:ILE:HD12	1.90	0.72
1:A:996:ASN:HD22	1:A:1021:MET:HB2	1.55	0.72
1:A:74:ARG:NH1	1:A:346:ILE:HD13	2.04	0.72
1:A:838:GLN:NE2	1:A:842:ASP:OD2	2.24	0.71
1:A:72:GLU:HG2	1:A:350:ARG:HD3	1.71	0.70
1:A:457:LYS:HD2	1:A:938:LYS:HB2	1.73	0.70
1:A:534:HIS:CE1	1:A:562:LEU:HG	2.26	0.70
1:A:574:THR:HA	1:A:588:ILE:H	1.56	0.70
1:A:219:ASP:O	1:A:222:LYS:HB3	1.92	0.69
1:A:233:THR:HA	1:A:235:ARG:HH12	1.56	0.69
1:A:481:LEU:HD22	1:A:890:THR:HG22	1.75	0.69
1:A:599:TYR:HE1	1:A:613:LEU:HB2	1.57	0.69
1:A:599:TYR:CE1	1:A:613:LEU:HB2	2.29	0.68
1:A:530:GLU:HB2	1:A:611:TYR:OH	1.92	0.68
1:A:22:LEU:HD11	1:A:125:GLY:HA3	1.75	0.68
1:A:600:TYR:CD2	1:A:603:ASP:HA	2.29	0.68
1:A:593:GLU:HB3	1:A:598:LYS:HG3	1.76	0.68
1:A:409:CYS:O	1:A:414:GLN:NE2	2.28	0.67
1:A:858:TYR:HA	1:A:878:PRO:HA	1.77	0.67
1:A:68:MET:HG2	1:A:133:ALA:HB1	1.75	0.66
1:A:578:TRP:CE3	1:A:583:VAL:HG22	2.30	0.66
1:A:903:ILE:HA	1:A:906:ILE:HD12	1.76	0.66
1:A:99:LEU:HB3	1:A:106:LEU:HD23	1.77	0.66
1:A:218:LEU:HD12	1:A:230:LEU:HD11	1.78	0.65
1:A:637:ASN:O	1:A:640:PRO:HD3	1.96	0.65
1:A:39:GLU:HG3	1:A:41:ARG:HH11	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1027:LEU:HB2	1:A:1131:PHE:HZ	1.60	0.64
1:A:621:HIS:CD2	1:A:630:ARG:HB3	2.32	0.64
1:A:579:ARG:N	1:A:582:ARG:O	2.25	0.64
1:A:74:ARG:HH12	1:A:346:ILE:CD1	2.11	0.64
1:A:901:GLN:HA	1:A:904:ARG:HD3	1.78	0.64
1:A:18:ILE:HD12	1:A:121:ASN:CG	2.23	0.64
1:A:74:ARG:NH1	1:A:346:ILE:CD1	2.60	0.63
1:A:407:GLU:O	1:A:408:HIS:ND1	2.32	0.63
1:A:1167:LEU:HD21	1:A:1180:LEU:HD13	1.81	0.63
1:A:1040:PRO:HG2	1:A:1128:ASN:HB2	1.81	0.63
1:A:223:LYS:HG2	1:A:224:ASP:H	1.64	0.63
1:A:221:PHE:CE1	1:A:249:PHE:CE1	2.84	0.62
1:A:999:PRO:HD3	1:A:1020:TYR:HB3	1.82	0.62
1:A:1027:LEU:HB2	1:A:1131:PHE:CZ	2.35	0.62
1:A:68:MET:CG	1:A:133:ALA:HB1	2.30	0.62
1:A:170:LEU:HD23	1:A:177:VAL:HG22	1.81	0.62
1:A:599:TYR:CZ	1:A:607:PHE:HB2	2.33	0.62
1:A:45:VAL:HG22	1:A:52:VAL:HG22	1.81	0.62
1:A:578:TRP:NE1	1:A:581:GLY:H	1.93	0.62
1:A:82:PHE:HE1	1:A:110:SER:HB3	1.65	0.61
1:A:588:ILE:HA	1:A:601:LEU:HD21	1.83	0.61
1:A:130:HIS:O	1:A:134:MET:HG2	2.00	0.61
1:A:221:PHE:CE1	1:A:249:PHE:CD1	2.89	0.61
1:A:493:ARG:HD2	1:A:509:ILE:HG21	1.82	0.61
1:A:596:THR:OG1	1:A:598:LYS:NZ	2.34	0.61
1:A:190:ILE:HG23	1:A:204:LEU:HD22	1.84	0.60
1:A:857:THR:HA	1:A:879:LYS:HD2	1.83	0.60
1:A:150:ILE:O	1:A:153:VAL:HG12	2.01	0.60
1:A:621:HIS:CD2	1:A:630:ARG:HG2	2.36	0.60
1:A:303:ASP:OD2	1:A:305:LYS:NZ	2.35	0.60
1:A:1089:ILE:HG21	1:A:1124:ILE:HD11	1.83	0.60
1:A:859:ASN:OD1	1:A:879:LYS:NZ	2.35	0.60
1:A:1022:GLN:OE1	1:A:1133:ARG:NH1	2.31	0.59
1:A:1023:MET:HE3	1:A:1131:PHE:HE2	1.67	0.59
1:A:1090:CYS:HB2	1:A:1131:PHE:HB2	1.84	0.59
1:A:286:PHE:O	1:A:290:PHE:N	2.34	0.59
1:A:621:HIS:CD2	1:A:630:ARG:N	2.70	0.59
1:A:707:SER:HB2	1:A:1080:ILE:HB	1.84	0.59
1:A:385:VAL:O	1:A:389:ILE:HG13	2.03	0.59
1:A:329:THR:HB	1:A:345:TYR:CE1	2.37	0.59
1:A:329:THR:HB	1:A:345:TYR:HE1	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:578:TRP:HA	1:A:583:VAL:HA	1.85	0.58
1:A:71:LYS:HB3	1:A:72:GLU:OE1	2.03	0.58
1:A:72:GLU:HG2	1:A:350:ARG:CD	2.33	0.58
1:A:221:PHE:HE1	1:A:249:PHE:CD1	2.21	0.58
1:A:530:GLU:HG3	1:A:531:LYS:N	2.19	0.58
1:A:641:HIS:CD2	1:A:713:LEU:HD12	2.38	0.58
1:A:211:PHE:HE2	1:A:236:PRO:HA	1.68	0.58
1:A:589:ARG:HB2	1:A:600:TYR:CE2	2.39	0.58
1:A:648:TYR:HE2	1:A:670:LEU:HD22	1.67	0.58
1:A:876:LEU:HD23	1:A:887:GLU:H	1.68	0.57
1:A:975:LEU:HD12	1:A:1050:PRO:HG3	1.86	0.57
1:A:363:GLY:H	1:A:413:GLN:HE22	1.52	0.57
1:A:170:LEU:HD12	1:A:173:ILE:HD11	1.86	0.57
1:A:186:LYS:O	1:A:190:ILE:HG12	2.03	0.57
1:A:872:PHE:HB3	1:A:896:LEU:HD22	1.87	0.57
1:A:456:LYS:HA	1:A:939:PRO:HG2	1.86	0.57
1:A:477:GLN:NE2	1:A:898:GLU:OE2	2.35	0.57
1:A:480:GLU:HB2	1:A:493:ARG:HH11	1.70	0.56
1:A:222:LYS:HE2	1:A:231:GLY:O	2.04	0.56
1:A:350:ARG:HA	1:A:396:THR:HG21	1.87	0.56
1:A:351:MET:SD	1:A:1014:PHE:CD2	2.99	0.56
1:A:496:CYS:HB3	1:A:503:LEU:HD21	1.87	0.56
1:A:530:GLU:HG2	1:A:532:TRP:NE1	2.20	0.56
1:A:749:ILE:O	1:A:753:TYR:OH	2.24	0.56
1:A:488:ASP:OD2	1:A:665:ARG:NH2	2.39	0.56
1:A:1066:LYS:NZ	1:A:1117:GLN:O	2.34	0.56
1:A:1067:VAL:HA	1:A:1182:VAL:HG12	1.87	0.56
1:A:1150:HIS:HD1	1:A:1152:THR:HG23	1.71	0.56
1:A:648:TYR:HB3	1:A:651:LEU:HG	1.88	0.56
1:A:671:ILE:HD12	1:A:682:ILE:HB	1.87	0.56
1:A:1050:PRO:C	1:A:1051:MET:HE2	2.32	0.55
1:A:1112:ILE:HG23	1:A:1112:ILE:O	2.06	0.55
1:A:296:SER:O	1:A:1164:SER:OG	2.24	0.55
1:A:530:GLU:CG	1:A:531:LYS:N	2.70	0.55
1:A:50:ARG:HH22	1:A:140:THR:HG22	1.71	0.55
1:A:580:SER:O	1:A:580:SER:OG	2.23	0.55
1:A:591:THR:HG23	1:A:600:TYR:OH	2.07	0.55
1:A:55:SER:HB2	1:A:62:GLU:H	1.71	0.55
1:A:71:LYS:CB	1:A:72:GLU:OE1	2.55	0.55
1:A:401:VAL:HG12	1:A:451:ILE:HG12	1.88	0.55
1:A:561:PHE:HB3	1:A:631:LEU:HD22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:LEU:HD13	1:A:601:LEU:HD11	1.89	0.55
1:A:579:ARG:HD2	1:A:584:GLN:HB2	1.89	0.55
1:A:1063:LEU:HD22	1:A:1065:VAL:HG23	1.89	0.55
1:A:221:PHE:CE2	1:A:290:PHE:HE2	2.25	0.55
1:A:429:LEU:HD21	1:A:933:LEU:HD11	1.87	0.55
1:A:530:GLU:CD	1:A:531:LYS:H	2.15	0.55
1:A:621:HIS:CD2	1:A:630:ARG:CB	2.89	0.55
1:A:530:GLU:CG	1:A:531:LYS:H	2.20	0.54
1:A:1001:ARG:NH2	1:A:1051:MET:SD	2.81	0.54
1:A:648:TYR:CE2	1:A:670:LEU:HD22	2.42	0.54
1:A:858:TYR:HD2	1:A:876:LEU:HB3	1.72	0.54
1:A:951:ASP:OD2	1:A:953:ARG:NH1	2.41	0.54
1:A:233:THR:HA	1:A:235:ARG:NH1	2.22	0.54
1:A:347:ARG:O	1:A:351:MET:HB2	2.08	0.54
1:A:940:THR:O	1:A:956:ARG:NH1	2.40	0.54
1:A:90:GLN:HG2	1:A:112:ALA:HB1	1.90	0.54
1:A:839:ILE:O	1:A:843:ASN:ND2	2.41	0.54
1:A:30:VAL:HG22	1:A:111:LEU:HD22	1.90	0.54
1:A:34:ARG:HG2	1:A:35:LYS:HG2	1.90	0.54
1:A:374:TRP:O	1:A:478:GLN:NE2	2.36	0.54
1:A:993:ASP:OD1	1:A:993:ASP:N	2.39	0.53
1:A:1023:MET:HE3	1:A:1131:PHE:CE2	2.43	0.53
1:A:632:THR:OG1	1:A:633:ASP:OD2	2.22	0.53
1:A:299:ASN:ND2	1:A:1162:PHE:O	2.38	0.53
1:A:571:ASN:O	1:A:571:ASN:ND2	2.41	0.53
1:A:572:ASP:OD2	1:A:587:ARG:NH1	2.42	0.53
1:A:82:PHE:CE1	1:A:110:SER:HB3	2.42	0.53
1:A:216:SER:HB3	1:A:1112:ILE:CD1	2.37	0.53
1:A:243:LEU:HA	1:A:269:MET:HE1	1.90	0.53
1:A:74:ARG:HD3	1:A:392:HIS:NE2	2.24	0.53
1:A:291:LEU:HD23	1:A:1071:ARG:HH21	1.75	0.53
1:A:863:ALA:H	1:A:874:PHE:HA	1.74	0.53
1:A:502:LYS:HE2	1:A:852:ILE:HD11	1.90	0.52
1:A:534:HIS:HE1	1:A:562:LEU:HG	1.72	0.52
1:A:1167:LEU:HB2	1:A:1178:ALA:HB1	1.92	0.52
1:A:214:GLN:HB2	1:A:217:ILE:HD11	1.92	0.52
1:A:938:LYS:O	1:A:955:ILE:N	2.41	0.52
1:A:146:LEU:HD11	1:A:206:TYR:HD2	1.73	0.52
1:A:605:LEU:HB2	1:A:617:TYR:OH	2.10	0.52
1:A:738:GLU:OE1	1:A:738:GLU:N	2.30	0.52
1:A:532:TRP:HB3	1:A:634:PRO:HB3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:ILE:HG21	1:A:278:ARG:HD2	1.92	0.51
1:A:164:ARG:HD2	1:A:554:THR:OG1	2.10	0.51
1:A:892:ARG:HH12	1:A:893:VAL:HG12	1.76	0.51
1:A:863:ALA:HB2	1:A:875:ILE:HG13	1.92	0.51
1:A:351:MET:SD	1:A:1014:PHE:HD2	2.34	0.51
1:A:500:ASP:O	1:A:502:LYS:HG3	2.11	0.51
1:A:860:VAL:HG11	1:A:903:ILE:HB	1.93	0.51
1:A:190:ILE:HG13	1:A:205:PHE:HB2	1.93	0.51
1:A:221:PHE:CE2	1:A:290:PHE:CE2	2.99	0.51
1:A:18:ILE:HD12	1:A:121:ASN:OD1	2.12	0.50
1:A:218:LEU:HD23	1:A:291:LEU:HD12	1.93	0.50
1:A:381:LYS:O	1:A:385:VAL:HG23	2.12	0.50
1:A:400:PRO:HG3	1:A:447:LEU:HB3	1.93	0.50
1:A:534:HIS:HB3	1:A:637:ASN:HD21	1.76	0.50
1:A:28:MET:HE3	1:A:113:ALA:HB2	1.92	0.50
1:A:206:TYR:CE1	1:A:210:MET:HE3	2.46	0.50
1:A:218:LEU:CD2	1:A:291:LEU:HD12	2.42	0.50
1:A:624:CYS:SG	1:A:625:ALA:N	2.83	0.50
1:A:18:ILE:CD1	1:A:121:ASN:CG	2.85	0.50
1:A:1061:MET:HE2	1:A:1127:PRO:HB3	1.94	0.50
1:A:350:ARG:HE	1:A:392:HIS:HD2	1.60	0.50
1:A:22:LEU:CD1	1:A:125:GLY:HA3	2.42	0.49
1:A:985:VAL:HG11	1:A:1002:LEU:HD21	1.94	0.49
1:A:840:ILE:HG23	1:A:846:GLY:C	2.37	0.49
1:A:71:LYS:O	1:A:72:GLU:CD	2.54	0.49
1:A:943:THR:HG22	1:A:946:ASN:HB3	1.94	0.49
1:A:872:PHE:CE1	1:A:893:VAL:HB	2.47	0.49
1:A:880:GLN:OE1	1:A:883:ASP:N	2.45	0.49
1:A:151:TYR:CZ	1:A:157:ARG:HB2	2.48	0.49
1:A:221:PHE:CZ	1:A:229:ILE:HG21	2.48	0.49
1:A:711:GLU:N	1:A:715:GLU:OE1	2.45	0.49
1:A:210:MET:O	1:A:214:GLN:HG2	2.13	0.49
1:A:751:SER:OG	1:A:752:LEU:N	2.42	0.48
1:A:71:LYS:C	1:A:72:GLU:OE1	2.55	0.48
1:A:214:GLN:HB2	1:A:217:ILE:CD1	2.43	0.48
1:A:621:HIS:NE2	1:A:630:ARG:CG	2.75	0.48
1:A:71:LYS:C	1:A:72:GLU:CD	2.80	0.48
1:A:483:MET:HE1	1:A:888:PHE:HE1	1.77	0.48
1:A:65:LEU:HD13	1:A:109:LEU:HD23	1.96	0.48
1:A:372:HIS:O	1:A:375:THR:OG1	2.31	0.48
1:A:947:LEU:HD12	1:A:977:TYR:CG	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1084:PHE:CE1	1:A:1137:TYR:HB2	2.48	0.48
1:A:617:TYR:CD2	1:A:622:LEU:HB3	2.49	0.48
1:A:203:HIS:ND1	1:A:207:LYS:HE3	2.28	0.48
1:A:613:LEU:HG	1:A:617:TYR:CE1	2.48	0.48
1:A:876:LEU:HD21	1:A:888:PHE:CE2	2.49	0.48
1:A:300:SER:HA	1:A:1164:SER:O	2.14	0.48
1:A:142:ILE:HG23	1:A:206:TYR:HE2	1.79	0.48
1:A:363:GLY:H	1:A:413:GLN:NE2	2.11	0.48
1:A:71:LYS:O	1:A:72:GLU:CG	2.62	0.47
1:A:507:ASP:OD1	1:A:508:ASP:N	2.38	0.47
1:A:577:PHE:HE1	1:A:586:CYS:HB2	1.79	0.47
1:A:673:LYS:C	1:A:673:LYS:HD3	2.39	0.47
1:A:855:LEU:HD23	1:A:906:ILE:HD13	1.96	0.47
1:A:892:ARG:HB2	1:A:895:GLU:HB2	1.96	0.47
1:A:214:GLN:HG3	1:A:291:LEU:HD11	1.97	0.47
1:A:324:SER:OG	1:A:353:CYS:HA	2.15	0.47
1:A:164:ARG:O	1:A:168:THR:HG23	2.14	0.47
1:A:544:GLU:OE2	1:A:545:LYS:NZ	2.47	0.47
1:A:1063:LEU:HD12	1:A:1155:ILE:HD11	1.95	0.47
1:A:66:ASP:OD1	1:A:148:LYS:NZ	2.41	0.47
1:A:358:LEU:HB2	1:A:405:ILE:HD13	1.95	0.47
1:A:663:ILE:HG22	1:A:743:TYR:CG	2.50	0.47
1:A:621:HIS:HD2	1:A:629:LEU:C	2.23	0.47
1:A:39:GLU:HB2	1:A:41:ARG:HD3	1.97	0.47
1:A:220:GLU:C	1:A:222:LYS:N	2.72	0.47
1:A:286:PHE:HB2	1:A:289:GLU:CD	2.40	0.47
1:A:541:THR:HA	1:A:544:GLU:HG2	1.97	0.47
1:A:876:LEU:HD11	1:A:888:PHE:HE2	1.80	0.46
1:A:127:LYS:O	1:A:131:GLN:HG2	2.15	0.46
1:A:405:ILE:HG21	1:A:417:MET:HE1	1.97	0.46
1:A:553:GLU:OE2	1:A:554:THR:HG22	2.16	0.46
1:A:559:GLY:HA2	1:A:578:TRP:O	2.15	0.46
1:A:46:ILE:O	1:A:50:ARG:N	2.48	0.46
1:A:1039:GLN:NE2	1:A:1043:MET:SD	2.86	0.46
1:A:17:GLN:HG3	1:A:20:ARG:HH12	1.81	0.46
1:A:351:MET:SD	1:A:1014:PHE:CE2	3.08	0.46
1:A:588:ILE:HA	1:A:601:LEU:CD2	2.44	0.46
1:A:836:GLU:HA	1:A:839:ILE:HD12	1.98	0.46
1:A:1085:VAL:HG22	1:A:1136:VAL:HG12	1.97	0.46
1:A:18:ILE:CD1	1:A:121:ASN:OD1	2.64	0.46
1:A:530:GLU:CG	1:A:532:TRP:CE2	2.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1019:LYS:HG3	1:A:1133:ARG:CZ	2.46	0.46
1:A:217:ILE:C	1:A:219:ASP:N	2.72	0.46
1:A:1014:PHE:HA	1:A:1021:MET:CE	2.45	0.46
1:A:164:ARG:NH2	1:A:165:GLU:OE2	2.48	0.46
1:A:958:PHE:HB2	1:A:985:VAL:HG22	1.97	0.46
1:A:728:LYS:HB2	1:A:728:LYS:HE2	1.66	0.46
1:A:71:LYS:O	1:A:72:GLU:HG3	2.16	0.45
1:A:74:ARG:HH12	1:A:346:ILE:HD12	1.81	0.45
1:A:1059:ILE:HG22	1:A:1125:TYR:HE1	1.82	0.45
1:A:1076:LEU:HD21	1:A:1148:LEU:HD23	1.97	0.45
1:A:364:PRO:HG3	1:A:379:LYS:NZ	2.32	0.45
1:A:1070:ALA:HB3	1:A:1111:PRO:HG2	1.98	0.45
1:A:146:LEU:HD11	1:A:206:TYR:CD2	2.51	0.45
1:A:903:ILE:O	1:A:907:THR:HG23	2.17	0.45
1:A:1082:CYS:SG	1:A:1106:ASP:N	2.84	0.45
1:A:560:THR:O	1:A:578:TRP:N	2.50	0.45
1:A:456:LYS:HD2	1:A:926:ILE:HD11	1.98	0.45
1:A:553:GLU:CD	1:A:700:ARG:HD3	2.42	0.45
1:A:1000:PHE:CE2	1:A:1091:GLY:HA2	2.51	0.45
1:A:561:PHE:CD1	1:A:631:LEU:HB3	2.51	0.45
1:A:621:HIS:CD2	1:A:630:ARG:CG	3.00	0.45
1:A:569:PHE:CD1	1:A:587:ARG:NH1	2.85	0.45
1:A:855:LEU:HD12	1:A:855:LEU:HA	1.83	0.45
1:A:1107:ASN:O	1:A:1111:PRO:HB3	2.17	0.45
1:A:599:TYR:OH	1:A:607:PHE:HB2	2.17	0.44
1:A:622:LEU:O	1:A:628:GLU:HA	2.16	0.44
1:A:301:ILE:HA	1:A:1163:ARG:HD2	1.99	0.44
1:A:589:ARG:HB2	1:A:600:TYR:CD2	2.51	0.44
1:A:1068:LEU:HD23	1:A:1068:LEU:HA	1.82	0.44
1:A:52:VAL:HB	1:A:65:LEU:HD22	1.98	0.44
1:A:630:ARG:NH1	1:A:631:LEU:O	2.51	0.44
1:A:367:LYS:HG3	1:A:416:HIS:NE2	2.32	0.44
1:A:941:SER:H	1:A:942:LYS:NZ	2.14	0.44
1:A:241:VAL:N	1:A:285:LEU:O	2.44	0.44
1:A:88:VAL:HG13	1:A:89:ARG:HD2	2.00	0.44
1:A:132:GLU:O	1:A:136:ALA:N	2.50	0.44
1:A:219:ASP:O	1:A:222:LYS:N	2.45	0.44
1:A:544:GLU:HB2	1:A:583:VAL:HG11	1.99	0.44
1:A:83:GLU:HB2	1:A:84:ARG:HH12	1.83	0.44
1:A:1162:PHE:HB3	1:A:1181:LEU:HD11	2.00	0.44
1:A:530:GLU:HG3	1:A:531:LYS:H	1.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1022:GLN:HG2	1:A:1150:HIS:CE1	2.52	0.44
1:A:174:ASN:OD1	1:A:1108:GLY:HA3	2.18	0.44
1:A:899:TRP:O	1:A:902:SER:OG	2.34	0.44
1:A:1014:PHE:HA	1:A:1021:MET:HE3	2.00	0.44
1:A:320:HIS:HB3	1:A:1044:ARG:HD3	1.99	0.43
1:A:345:TYR:OH	1:A:357:GLU:O	2.25	0.43
1:A:1136:VAL:O	1:A:1148:LEU:HB2	2.18	0.43
1:A:386:VAL:HG13	1:A:424:VAL:HG21	2.00	0.43
1:A:454:LYS:HE2	1:A:955:ILE:HG21	2.01	0.43
1:A:532:TRP:HZ2	1:A:611:TYR:CE2	2.36	0.43
1:A:1059:ILE:HG13	1:A:1189:VAL:HG21	1.98	0.43
1:A:562:LEU:C	1:A:562:LEU:HD23	2.43	0.43
1:A:721:GLU:OE1	1:A:722:LYS:HG2	2.18	0.43
1:A:1018:ASP:OD1	1:A:1019:LYS:N	2.46	0.43
1:A:1138:GLU:HB3	1:A:1148:LEU:HD11	2.00	0.43
1:A:181:LYS:HE3	1:A:181:LYS:HB2	1.74	0.43
1:A:532:TRP:CH2	1:A:610:ILE:HB	2.53	0.43
1:A:455:HIS:HE1	1:A:926:ILE:HG23	1.83	0.43
1:A:447:LEU:HA	1:A:447:LEU:HD23	1.81	0.43
1:A:587:ARG:NE	1:A:589:ARG:HD3	2.27	0.43
1:A:945:ASP:OD1	1:A:945:ASP:C	2.60	0.43
1:A:163:LEU:HD11	1:A:167:LYS:HE3	2.00	0.43
1:A:545:LYS:HA	1:A:545:LYS:HD3	1.90	0.43
1:A:1023:MET:SD	1:A:1090:CYS:HB3	2.59	0.43
1:A:153:VAL:CG2	1:A:165:GLU:HB3	2.40	0.43
1:A:156:THR:O	1:A:157:ARG:HG2	2.19	0.43
1:A:221:PHE:CZ	1:A:249:PHE:CD1	3.07	0.43
1:A:491:TRP:CG	1:A:870:LYS:HZ3	2.37	0.43
1:A:941:SER:O	1:A:941:SER:OG	2.36	0.43
1:A:507:ASP:HB2	1:A:850:ARG:HG3	2.01	0.43
1:A:476:LYS:HE2	1:A:476:LYS:HB2	1.86	0.43
1:A:530:GLU:HG2	1:A:532:TRP:CZ2	2.54	0.43
1:A:532:TRP:HZ2	1:A:611:TYR:CE1	2.36	0.43
1:A:639:ASN:HB3	1:A:641:HIS:CE1	2.54	0.43
1:A:736:THR:O	1:A:740:LEU:N	2.47	0.43
1:A:275:ASP:HB3	1:A:278:ARG:HG3	2.01	0.42
1:A:696:ASN:HB2	1:A:703:VAL:HG12	1.99	0.42
1:A:669:PHE:HA	1:A:683:THR:O	2.19	0.42
1:A:707:SER:OG	1:A:1080:ILE:N	2.50	0.42
1:A:928:ILE:HD12	1:A:928:ILE:H	1.84	0.42
1:A:481:LEU:HD11	1:A:899:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:968:ARG:HA	1:A:968:ARG:HD3	1.82	0.42
1:A:1049:ASP:OD1	1:A:1049:ASP:N	2.53	0.42
1:A:1051:MET:HE2	1:A:1051:MET:N	2.34	0.42
1:A:980:LYS:HA	1:A:980:LYS:HD3	1.82	0.42
1:A:1005:CYS:SG	1:A:1051:MET:HE3	2.60	0.42
1:A:239:SER:O	1:A:287:VAL:N	2.50	0.42
1:A:618:ARG:HG2	1:A:631:LEU:O	2.18	0.42
1:A:1019:LYS:HG3	1:A:1133:ARG:NE	2.34	0.42
1:A:218:LEU:HD12	1:A:230:LEU:CD1	2.47	0.42
1:A:650:SER:O	1:A:651:LEU:HD23	2.19	0.42
1:A:1064:THR:OG1	1:A:1185:GLU:HB3	2.19	0.42
1:A:48:GLU:HB3	1:A:147:ARG:HG2	2.02	0.42
1:A:217:ILE:O	1:A:218:LEU:C	2.63	0.42
1:A:950:PRO:HB3	1:A:977:TYR:CD1	2.55	0.42
1:A:347:ARG:HH22	1:A:1173:GLU:CD	2.28	0.42
1:A:193:HIS:ND1	1:A:194:LYS:N	2.68	0.41
1:A:964:ASP:O	1:A:967:ILE:HG22	2.20	0.41
1:A:138:THR:OG1	1:A:296:SER:HB2	2.21	0.41
1:A:229:ILE:HG21	1:A:249:PHE:HB2	2.01	0.41
1:A:316:ASN:CG	1:A:1044:ARG:HH12	2.23	0.41
1:A:1062:THR:HG23	1:A:1187:ARG:HB2	2.02	0.41
1:A:221:PHE:CZ	1:A:290:PHE:HZ	2.38	0.41
1:A:327:HIS:CG	1:A:987:PRO:HD2	2.54	0.41
1:A:898:GLU:HA	1:A:901:GLN:OE1	2.20	0.41
1:A:1083:PRO:HA	1:A:1137:TYR:O	2.20	0.41
1:A:1089:ILE:HD11	1:A:1122:PHE:CD2	2.55	0.41
1:A:212:GLU:OE2	1:A:236:PRO:HB3	2.20	0.41
1:A:1017:ALA:HB2	1:A:1171:TYR:CD1	2.55	0.41
1:A:456:LYS:O	1:A:926:ILE:HD13	2.20	0.41
1:A:730:ARG:HD2	1:A:732:ARG:NH2	2.35	0.41
1:A:1087:VAL:HG22	1:A:1134:PHE:HD1	1.86	0.41
1:A:1138:GLU:OE1	1:A:1146:ASN:HB3	2.21	0.41
1:A:54:TRP:CE2	1:A:63:GLY:HA3	2.56	0.41
1:A:193:HIS:ND1	1:A:194:LYS:HG3	2.36	0.41
1:A:218:LEU:HD21	1:A:291:LEU:HB2	2.02	0.41
1:A:477:GLN:OE1	1:A:478:GLN:N	2.54	0.41
1:A:870:LYS:HD3	1:A:870:LYS:HA	1.86	0.41
1:A:256:GLU:OE1	1:A:258:TRP:NE1	2.51	0.41
1:A:335:GLN:HB3	1:A:372:HIS:CE1	2.56	0.41
1:A:535:LYS:HB3	1:A:536:LYS:H	1.65	0.41
1:A:610:ILE:H	1:A:610:ILE:HG12	1.72	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:876:LEU:HD13	1:A:876:LEU:HA	1.91	0.41
1:A:82:PHE:CZ	1:A:95:CYS:HB3	2.56	0.40
1:A:294:LEU:O	1:A:1164:SER:HB3	2.20	0.40
1:A:555:GLY:O	1:A:557:LYS:HG2	2.22	0.40
1:A:569:PHE:CG	1:A:587:ARG:HD3	2.56	0.40
1:A:160:SER:HA	1:A:197:LEU:O	2.21	0.40
1:A:1124:ILE:HD13	1:A:1124:ILE:HA	1.93	0.40
1:A:1137:TYR:HD1	1:A:1146:ASN:O	2.04	0.40
1:A:1137:TYR:CD1	1:A:1147:PHE:HA	2.55	0.40
1:A:171:PRO:HB3	1:A:1080:ILE:HD11	2.04	0.40
1:A:574:THR:HA	1:A:587:ARG:HA	2.03	0.40
1:A:663:ILE:HG22	1:A:743:TYR:CD1	2.57	0.40
1:A:212:GLU:HA	1:A:212:GLU:OE1	2.21	0.40
1:A:56:LYS:HA	1:A:56:LYS:HE2	2.04	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	1051/1265 (83%)	997 (95%)	54 (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	965/1158 (83%)	965 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	159	ASN
1	A	247	GLN
1	A	254	GLN
1	A	260	GLN
1	A	263	ASN
1	A	494	HIS
1	A	621	HIS
1	A	637	ASN
1	A	843	ASN
1	A	865	GLN
1	A	1025	HIS
1	A	1031	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

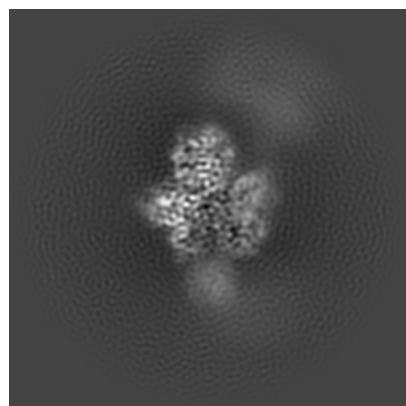
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-36571. These allow visual inspection of the internal detail of the map and identification of artifacts.

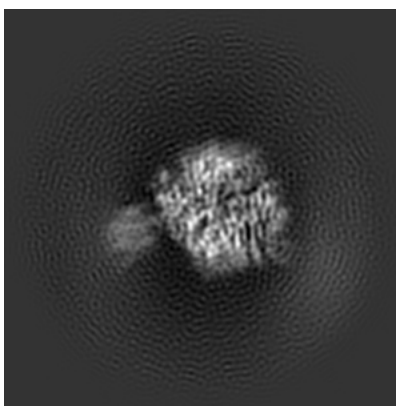
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

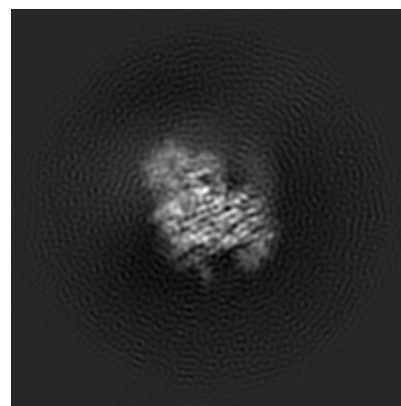
6.1.1 Primary map



X

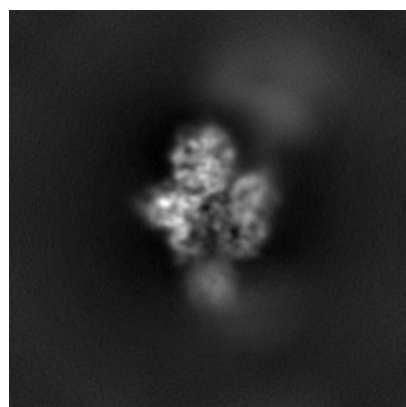


Y

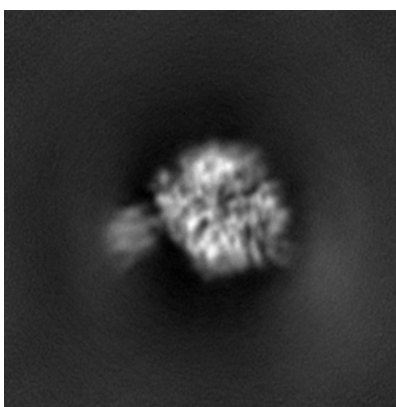


Z

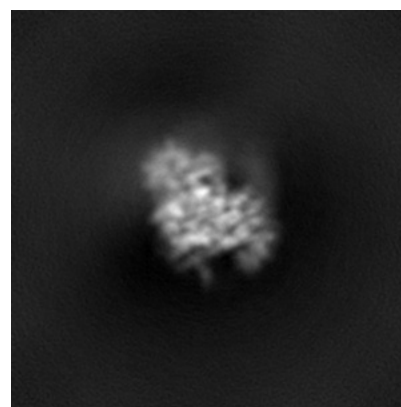
6.1.2 Raw map



X



Y

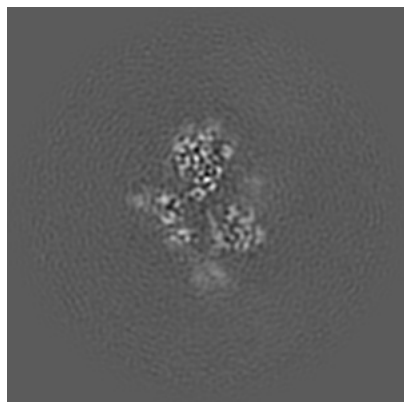


Z

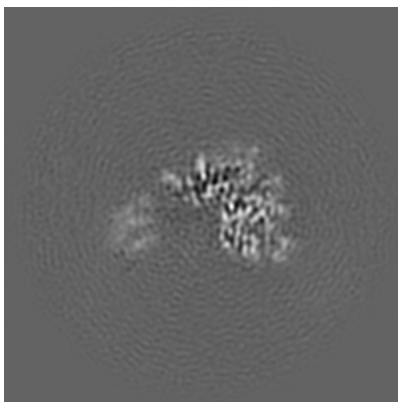
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

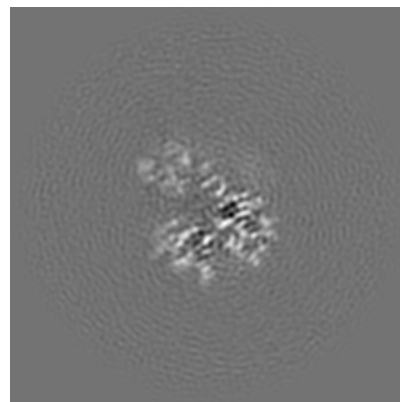
6.2.1 Primary map



X Index: 108

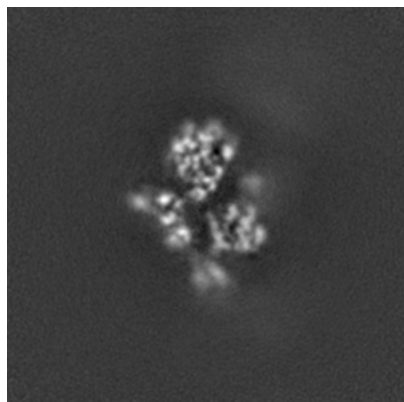


Y Index: 108

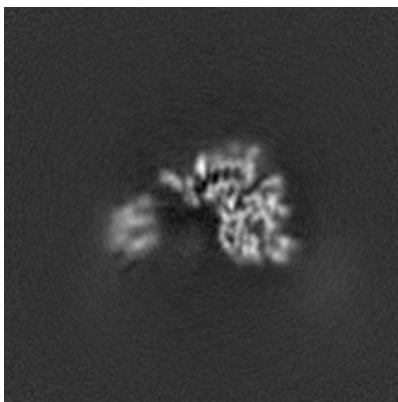


Z Index: 108

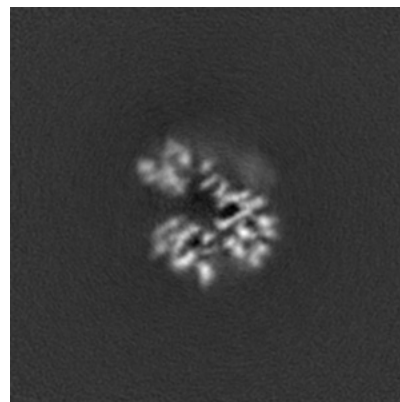
6.2.2 Raw map



X Index: 108



Y Index: 108

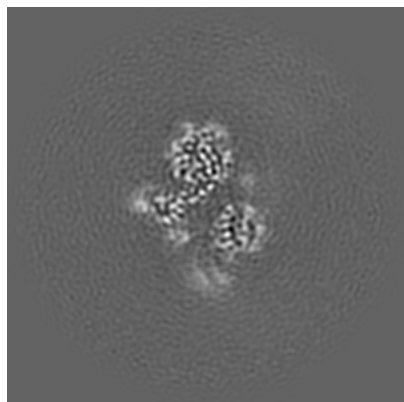


Z Index: 108

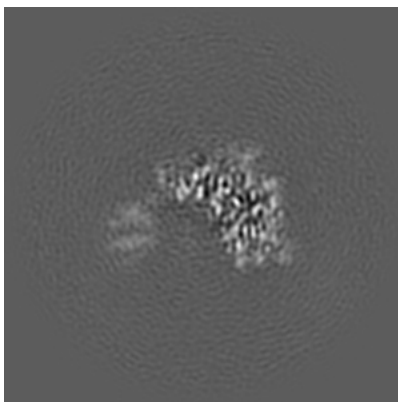
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

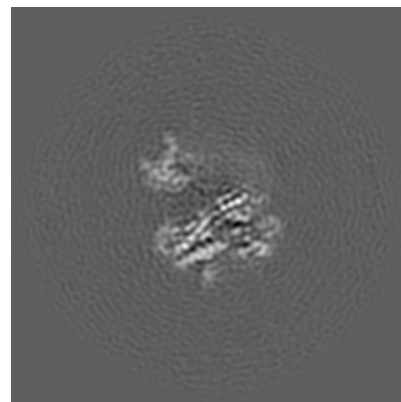
6.3.1 Primary map



X Index: 105

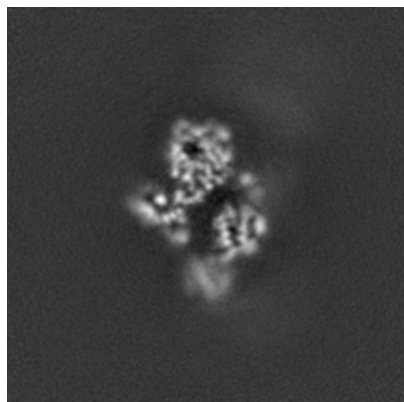


Y Index: 105

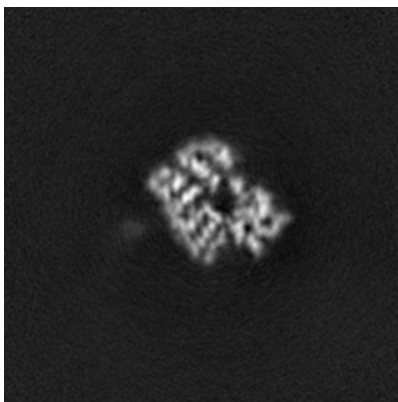


Z Index: 111

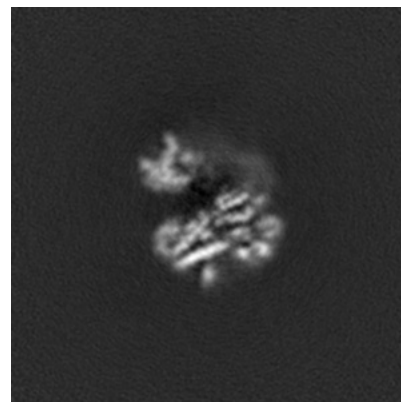
6.3.2 Raw map



X Index: 103



Y Index: 92

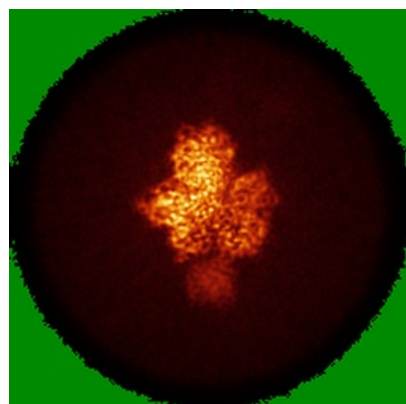


Z Index: 111

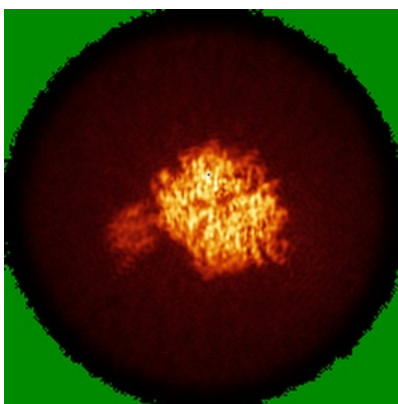
The images above show the largest variance slices of the map in three orthogonal directions.

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

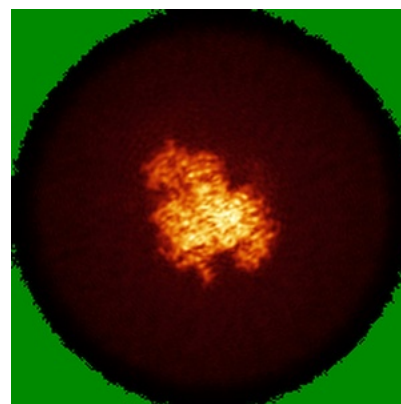
6.4.1 Primary map



X

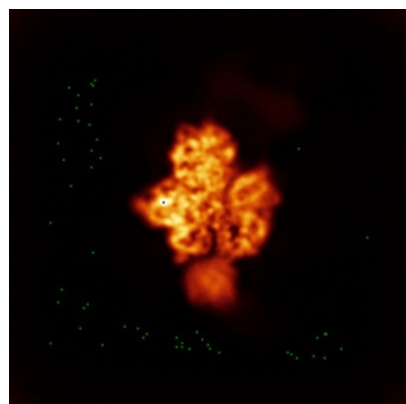


Y

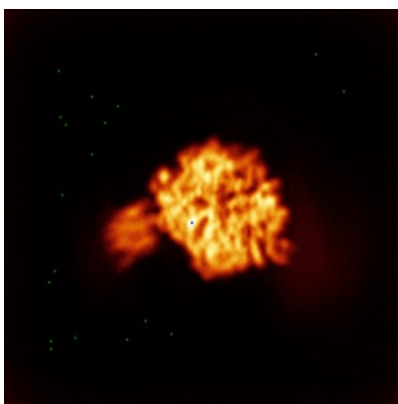


Z

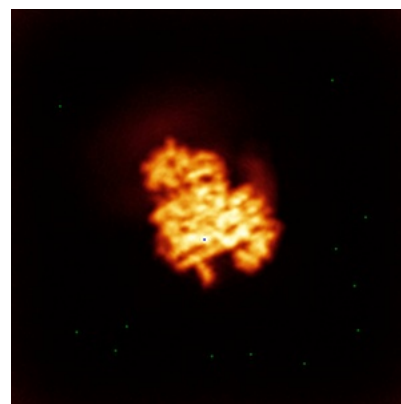
6.4.2 Raw map



X



Y



Z

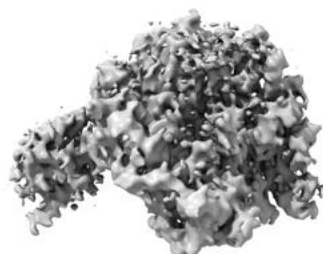
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

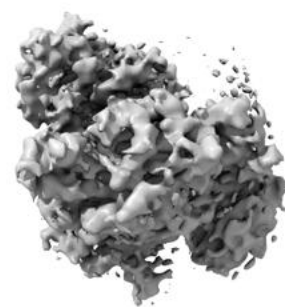
6.5.1 Primary map



X



Y



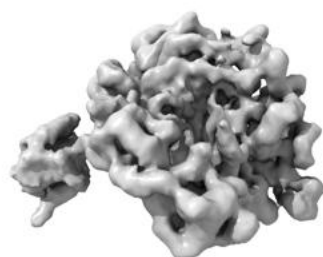
Z

The images above show the 3D surface view of the map at the recommended contour level 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

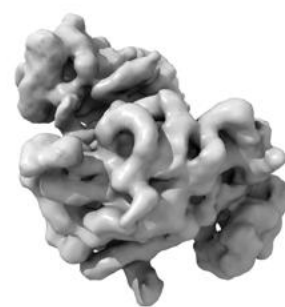
6.5.2 Raw map



X



Y



Z

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.

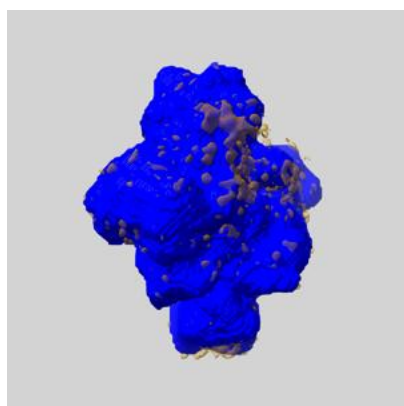
6.6 Mask visualisation [i](#)

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

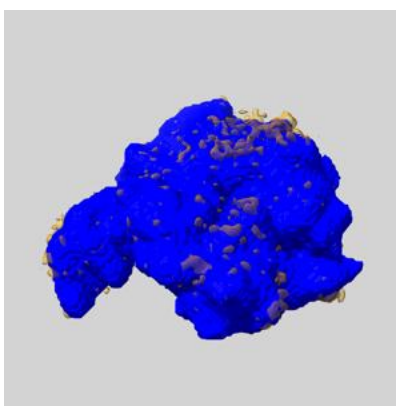
A mask typically either:

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

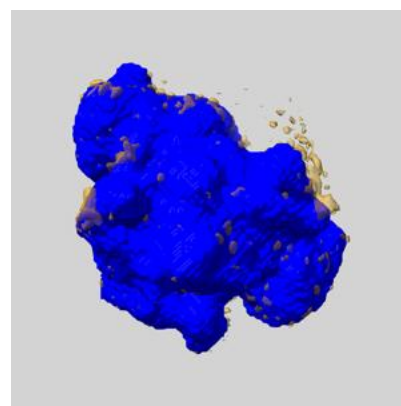
6.6.1 emd_36571_msk_1.map [i](#)



X



Y

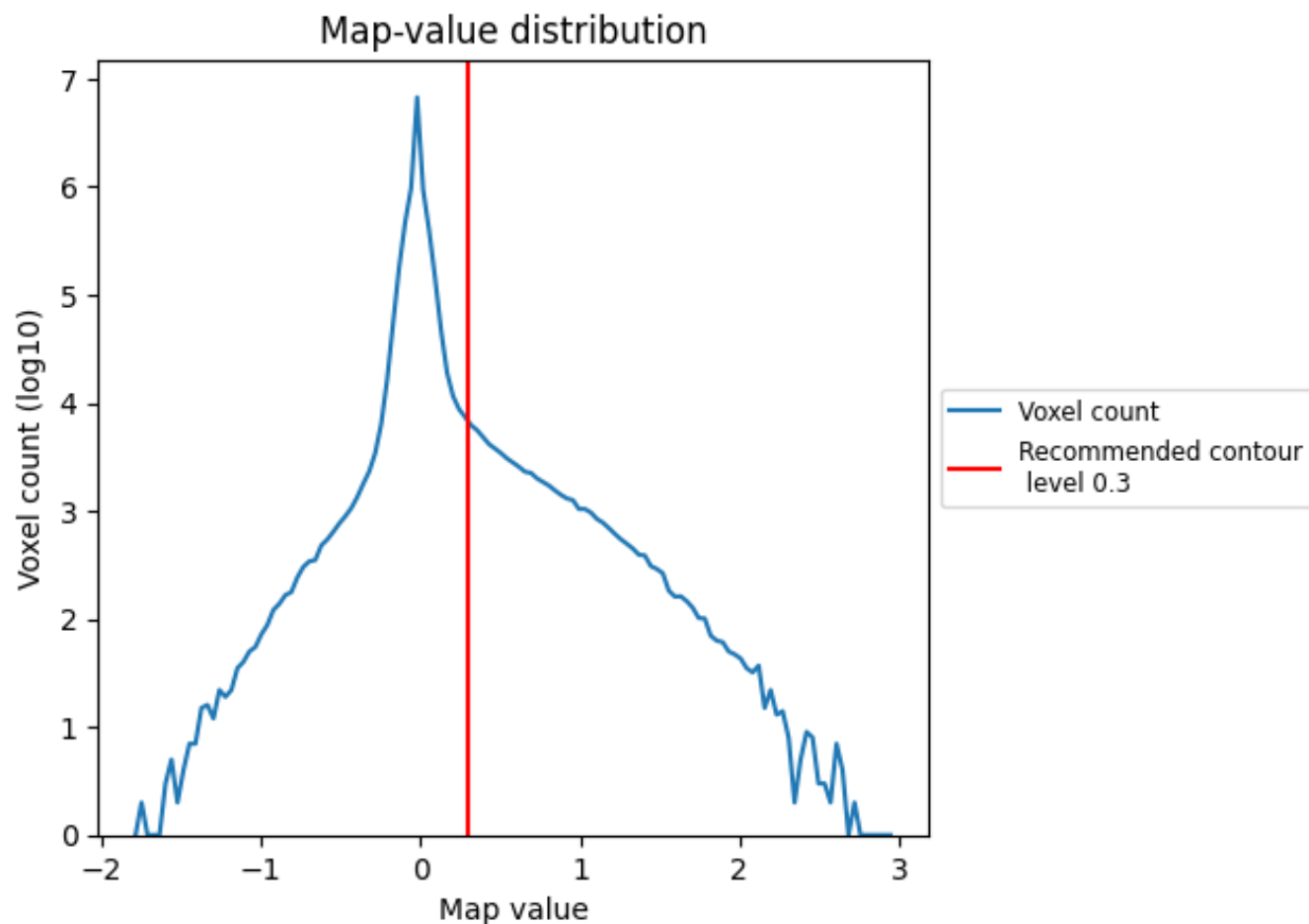


Z

7 Map analysis [i](#)

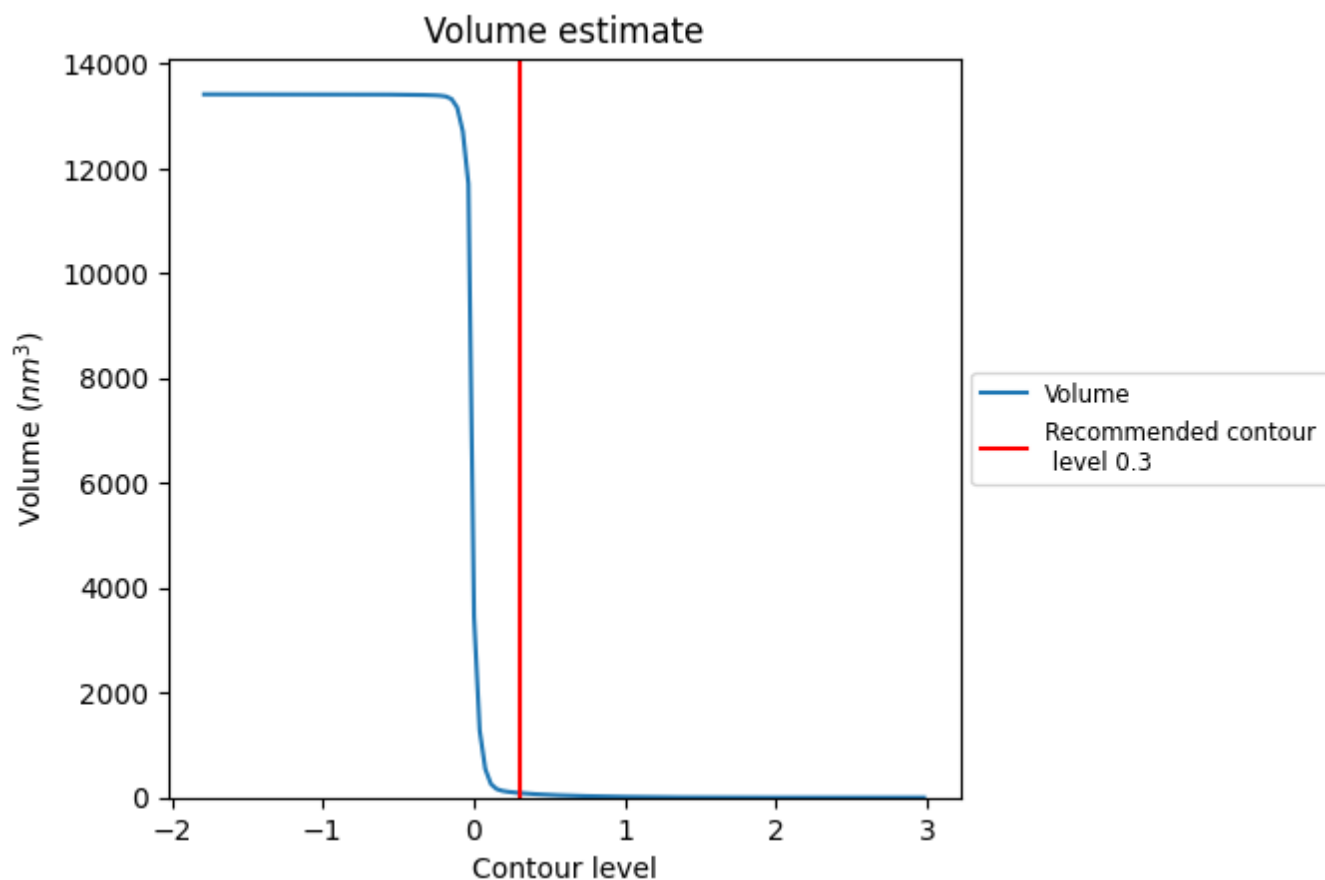
This section contains the results of statistical analysis of the map.

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.

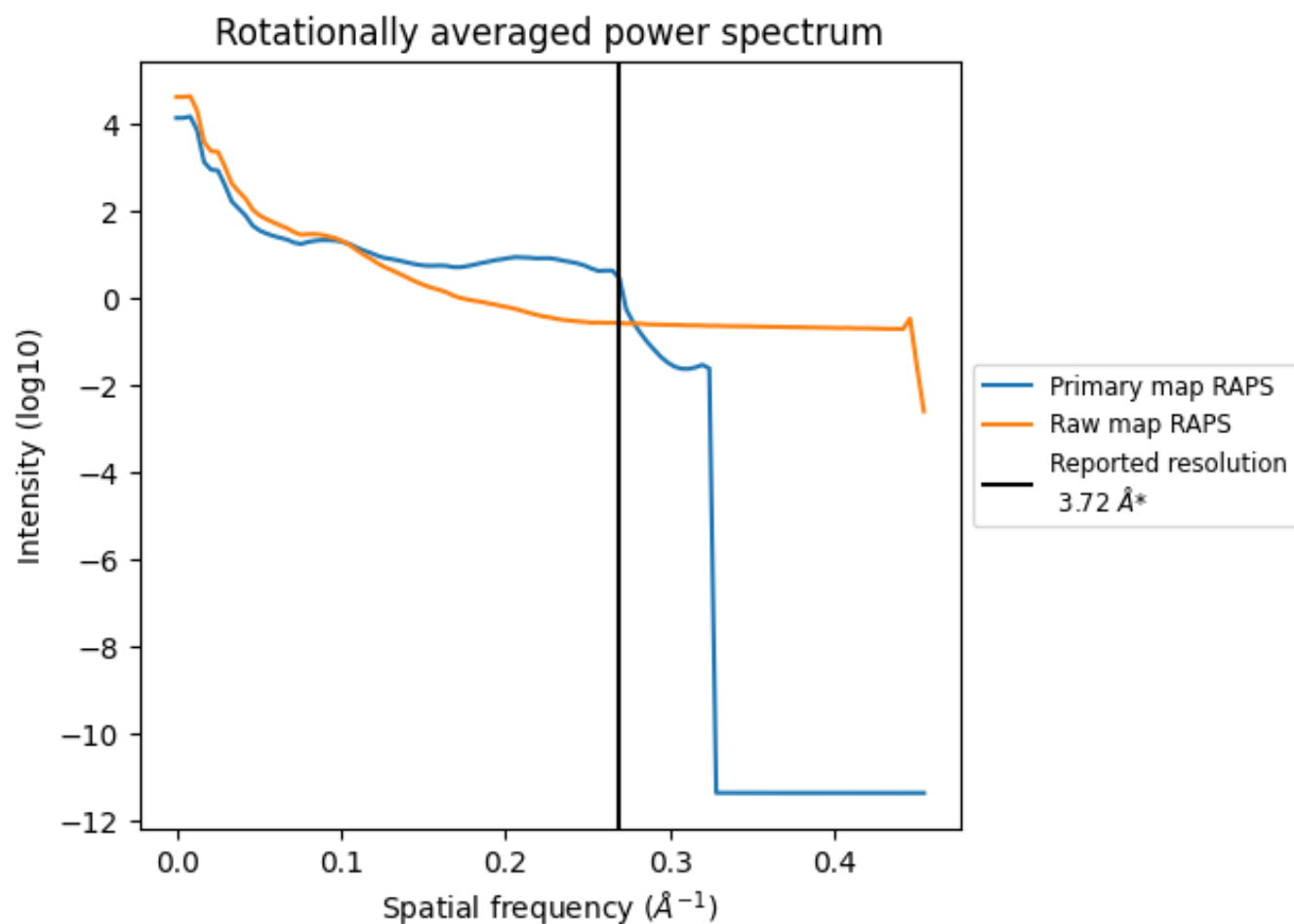
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 88 nm³; this corresponds to an approximate mass of 79 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 ⓘ

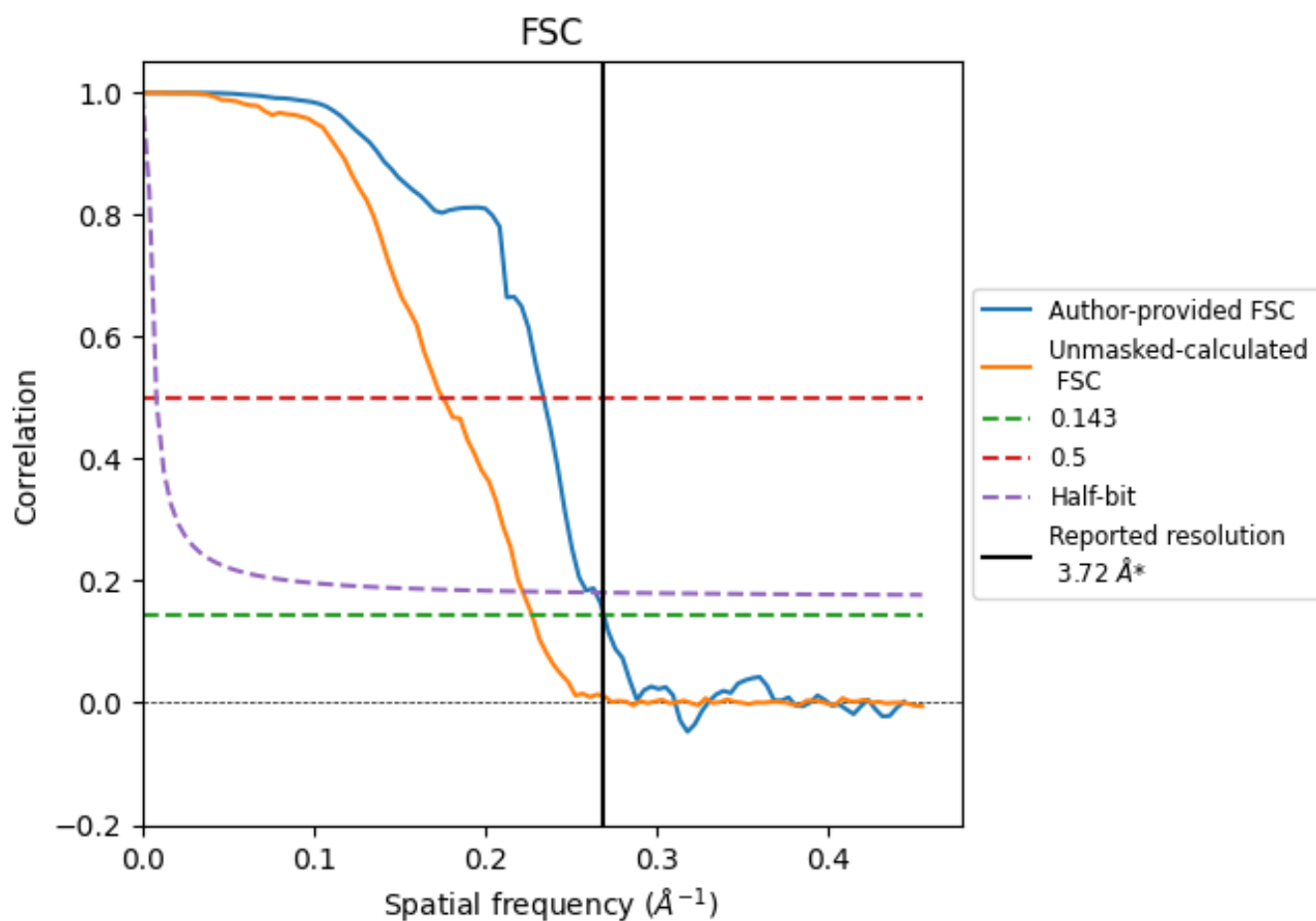


*Reported resolution corresponds to spatial frequency of 0.269 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.269 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

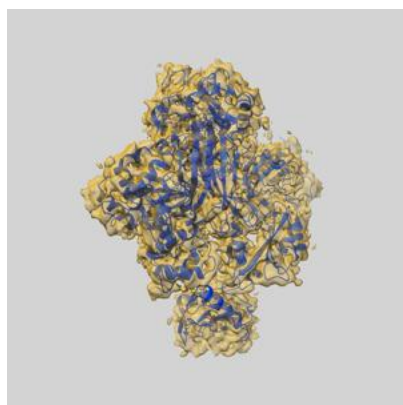
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.72	-	-
Author-provided FSC curve	3.72	4.28	3.79
Unmasked-calculated*	4.40	5.71	4.51

*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 4.40 differs from the reported value 3.72 by more than 10 %

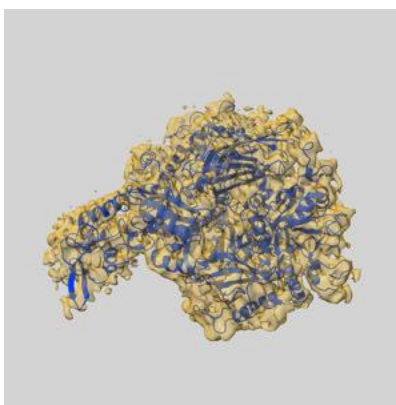
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-36571 and PDB model 8JQG. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

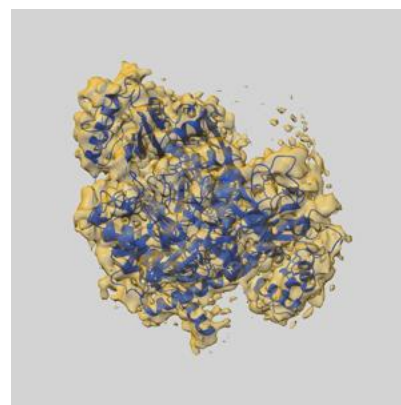
9.1 Map-model overlay [i](#)



X



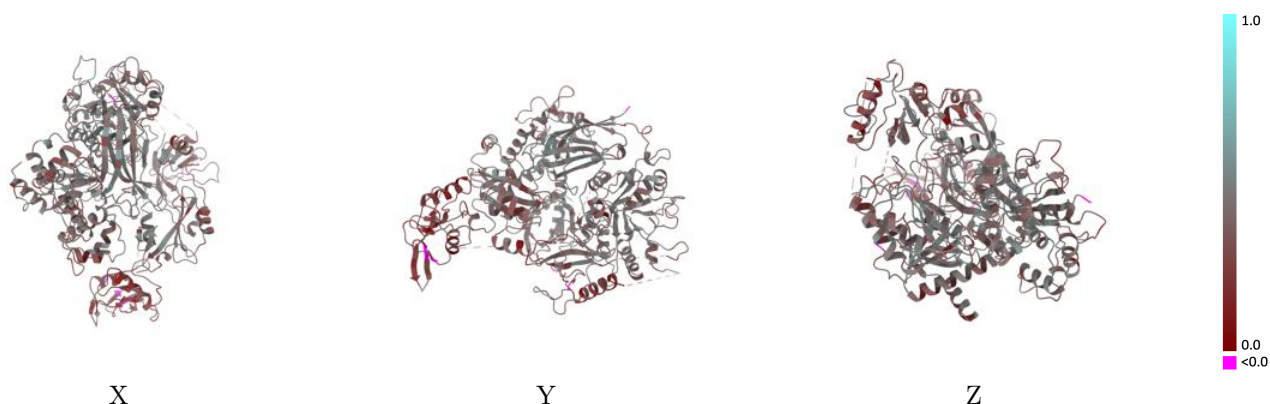
Y



Z

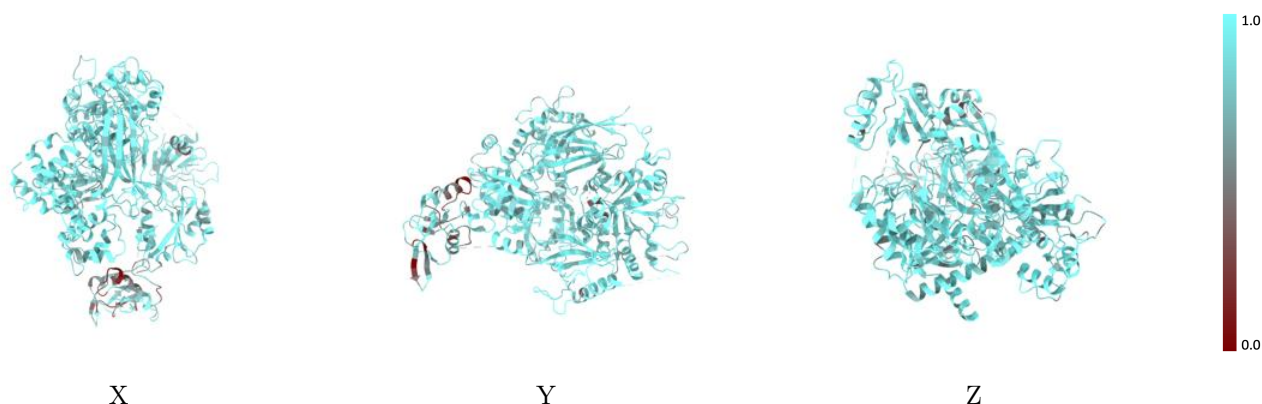
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



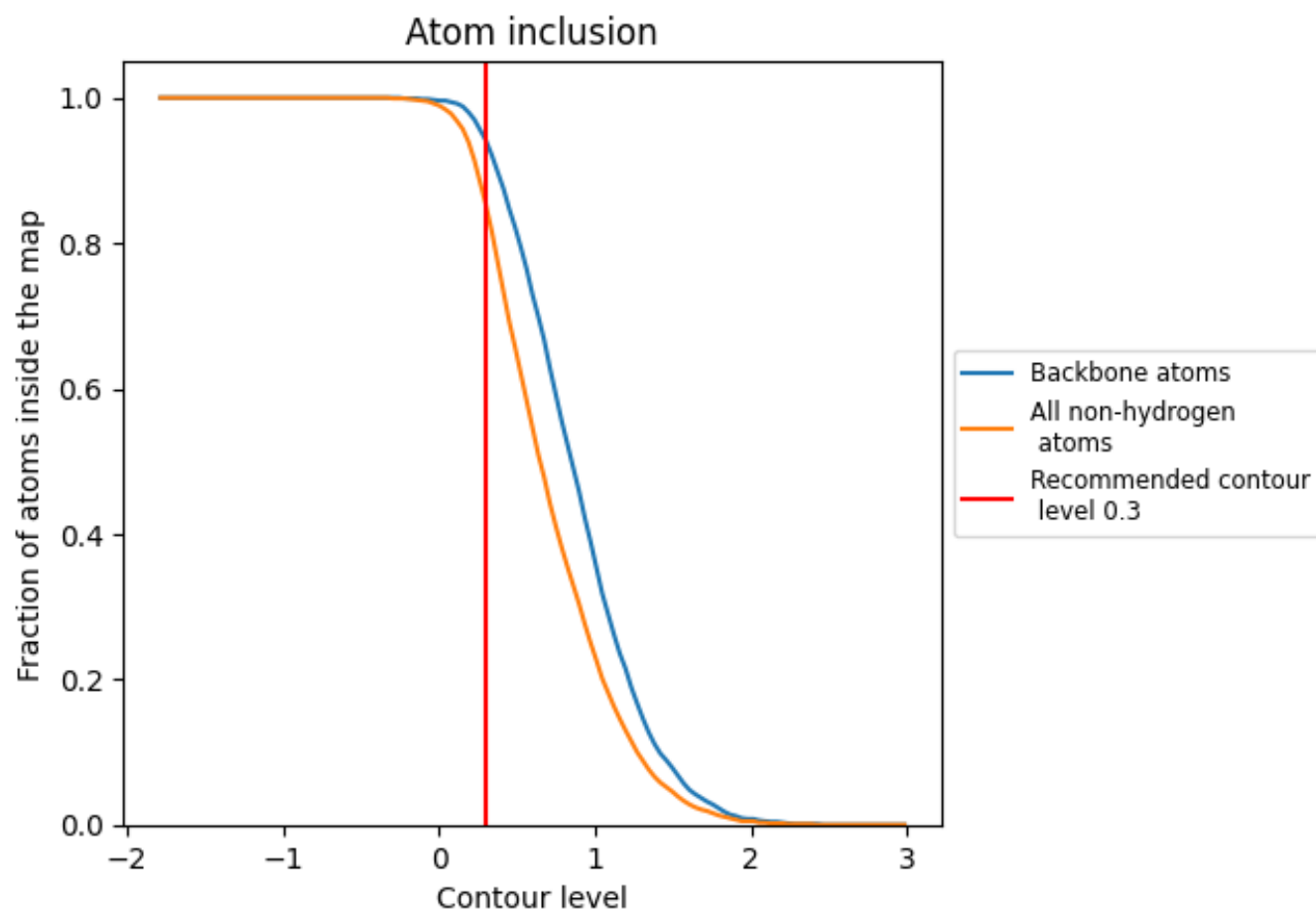
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 85% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8520	0.3840
A	0.8520	0.3840

