Full wwPDB EM Validation Report

Nov 9, 2026 – 10:20 AM EST

PDB ID	:	37MQ / pdb_000037mq
EMDB ID	:	EMD-78314
Title	:	Cryo-EM model of the minor reactivation conformation of Cob(II)-bound Thermus filiformis cobalamin-dependent methionine synthase (MetH)
Authors	:	Suksmith, P.; Wang, H.; Ando, N.
Deposited on	:	2026-07-27
Resolution	:	4.60 Å(reported)
Based on initial model	:	.

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

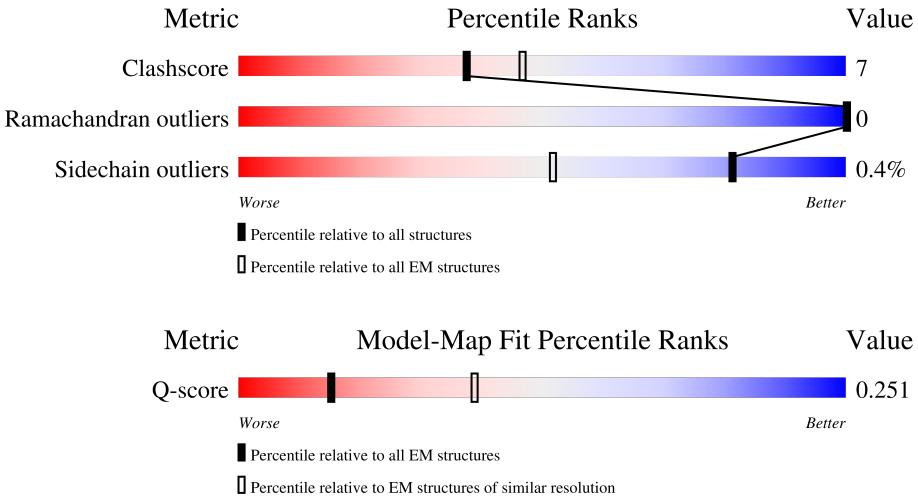
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.2
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data 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)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.60 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2407 (4.10 - 5.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	1178	20%, 81%, 14%, .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8989 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.

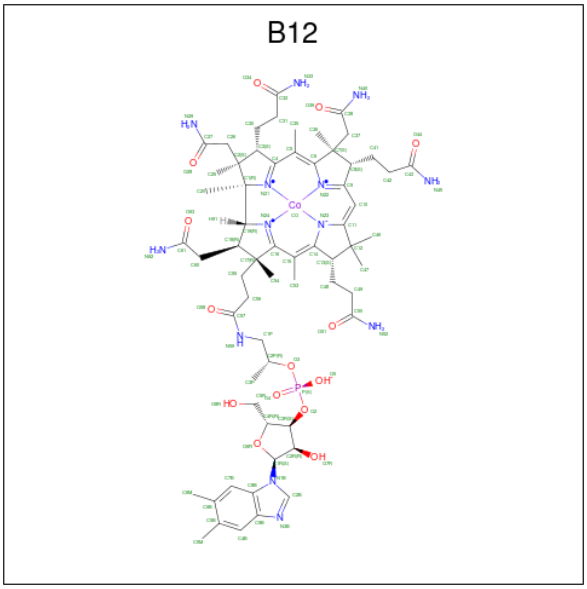
- Molecule 1 is a protein called Methionine synthase.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1133	8897	5663	1555	1646	33	0	0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
			Total	Zn	
2	A	1	1	1	0

- Molecule 3 is COBALAMIN (CCD ID: B12) (formula: $C_{62}H_{89}CoN_{13}O_{14}P$) (labeled as "Ligand of Interest" by depositor).



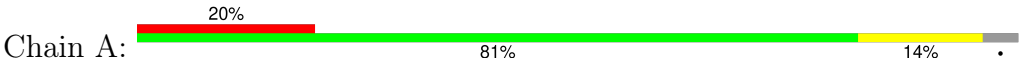
Mol	Chain	Residues	Atoms						AltConf
			Total	C	Co	N	O	P	
3	A	1	91	62	1	13	14	1	0

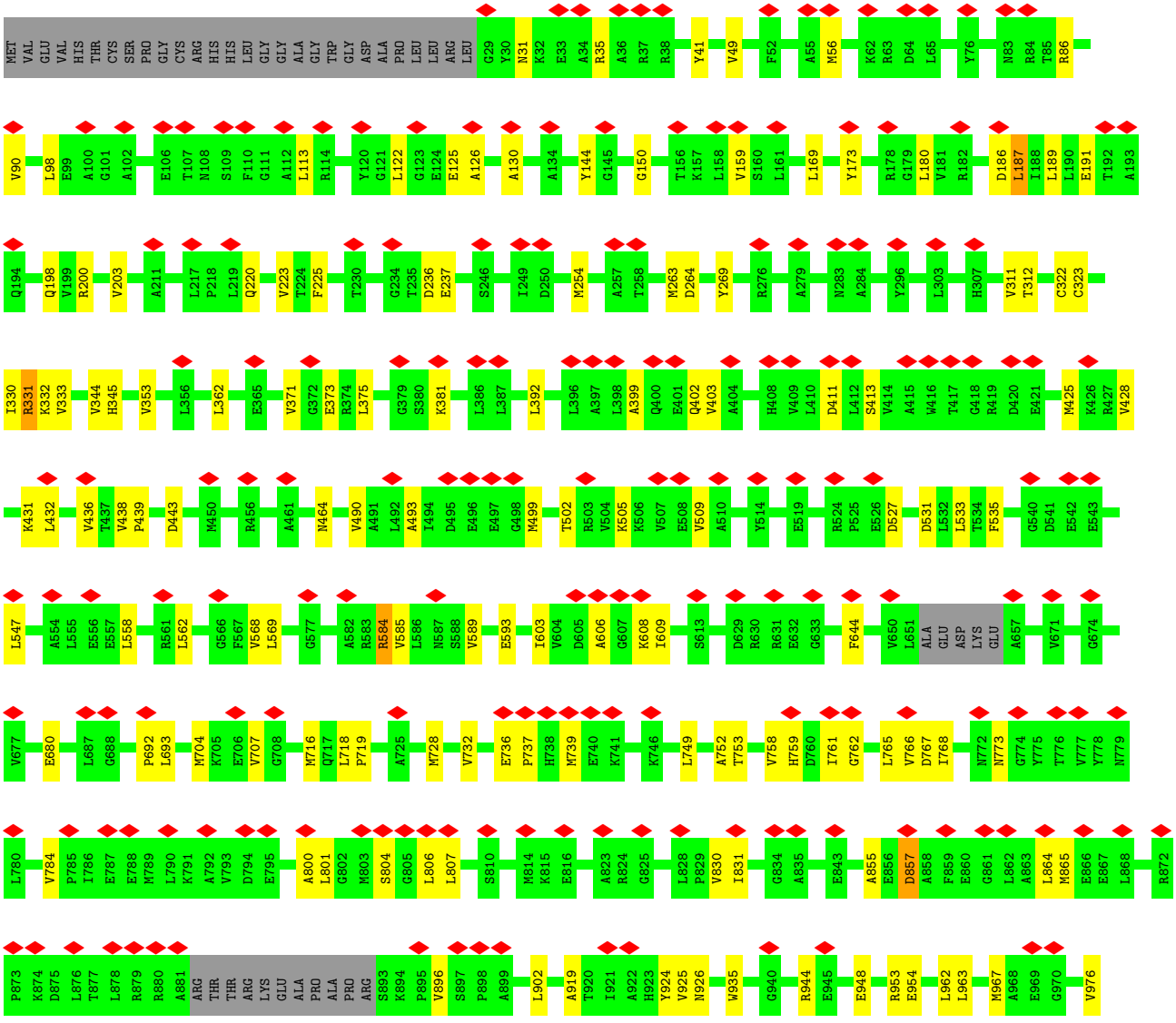
3 Residue-property plots (i)

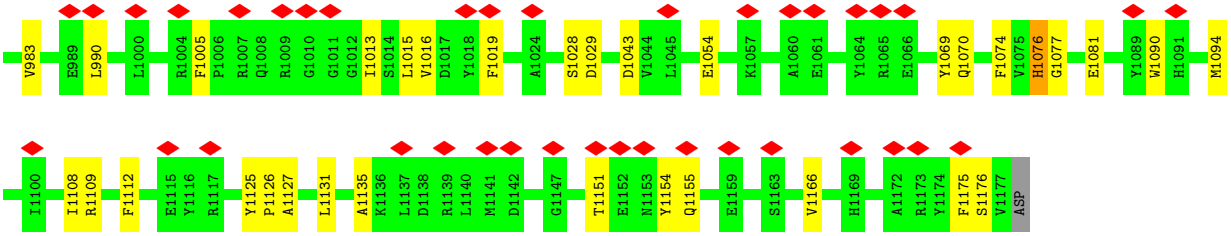
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

Chain A:






V983
E989
L990
L1000
R1004
F1005
P1006
R1007
Q1008
R1009
G1010
G1011
G1012
I1013
S1014
L1015
V1016
D1017
Y1018
F1019
A1024
S1028
D1029
D1043
V1044
L1045
E1054
K1057
A1060
E1061
Y1064
R1065
E1066
Y1069
Q1070
F1074
V1075
H1076
G1077
E1081
Y1089
W1090
H1091
M1094
I1100
I1108
R1109
F1112
E1115
Y1116
R1117
Y1125
P1126
A1127
L1131
A1135
K1136
L1137
D1138
R1139
L1140
M1141
D1142
G1147
T1151
E1152
N1153
Y1154
Q1155
E1159
S1163
V1166
H1169
A1172
R1173
Y1174
F1175
S1176
V1177
ASP

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	22471	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 ($e^-/Å^2$)	61.7	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.518	Depositor
Minimum map value	-0.123	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.018	Depositor
Recommended contour level	0.119	Depositor
Map size (Å)	273.92, 273.92, 273.92	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, B12

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	0/9077	0.52	0/12282

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	331	ARG	Sidechain
1	A	584	ARG	Sidechain
1	A	86	ARG	Sidechain

5.2 Too-close contacts

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	8897	0	8921	131	0
2	A	1	0	0	0	0
3	A	91	0	88	13	0
All	All	8989	0	9009	134	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 7.

All (134) 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:693:LEU:HG	1:A:739:MET:HE3	1.29	1.13
1:A:749:LEU:HD12	1:A:800:ALA:O	1.64	0.98
1:A:1126:PRO:HB2	1:A:1176:SER:HA	1.68	0.74
1:A:693:LEU:CG	1:A:739:MET:HE3	2.15	0.71
1:A:159:VAL:HG12	1:A:198:GLN:HG3	1.73	0.70
1:A:533:LEU:HD13	1:A:535:PHE:CZ	2.33	0.64
1:A:344:VAL:HG12	1:A:345:HIS:N	2.12	0.64
1:A:749:LEU:HD12	1:A:800:ALA:C	2.24	0.63
1:A:693:LEU:HG	1:A:739:MET:CE	2.19	0.62
1:A:189:LEU:HD11	1:A:191:GLU:HB3	1.81	0.61
1:A:983:VAL:HB	1:A:990:LEU:HD11	1.83	0.61
1:A:1125:TYR:HB3	1:A:1126:PRO:HD2	1.81	0.61
1:A:962:LEU:HD11	1:A:1074:PHE:HB3	1.83	0.59
1:A:1015:LEU:HD21	1:A:1127:ALA:HB1	1.84	0.59
1:A:704:MET:HE2	1:A:773:ASN:HD21	1.68	0.58
1:A:680:GLU:H	1:A:680:GLU:CD	2.12	0.57
1:A:831:ILE:HD11	1:A:855:ALA:HB2	1.84	0.57
1:A:896:VAL:HG23	1:A:1013:ILE:HD12	1.87	0.57
1:A:263:MET:HA	1:A:263:MET:HE2	1.87	0.56
1:A:1125:TYR:CE1	3:A:1202:B12:H352	2.41	0.56
1:A:56:MET:HE1	1:A:90:VAL:HG13	1.86	0.56
1:A:493:ALA:HA	1:A:509:VAL:HG11	1.89	0.55
1:A:150:GLY:HA3	1:A:180:LEU:HD23	1.88	0.55
1:A:759:HIS:HD2	3:A:1202:B12:H482	1.72	0.54
1:A:752:ALA:HB1	1:A:784:VAL:HG21	1.89	0.54
1:A:1155:GLN:HA	3:A:1202:B12:H402	1.73	0.54
1:A:990:LEU:HD13	1:A:1019:PHE:CD2	2.43	0.54
1:A:759:HIS:CE1	3:A:1202:B12:H412	2.43	0.53
1:A:413:SER:HB2	1:A:443:ASP:HB3	1.90	0.52
1:A:749:LEU:CD1	1:A:800:ALA:C	2.82	0.52
1:A:1094:MET:HE1	1:A:1166:VAL:HG21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:VAL:HG22	1:A:432:LEU:HD12	1.92	0.51
1:A:759:HIS:CD2	3:A:1202:B12:H482	2.45	0.51
1:A:353:VAL:HG23	1:A:439:PRO:HG2	1.93	0.51
1:A:392:LEU:HD12	1:A:431:LYS:CB	2.40	0.51
1:A:589:VAL:O	1:A:593:GLU:HG2	2.10	0.51
1:A:692:PRO:HB2	1:A:739:MET:HG2	1.92	0.51
1:A:225:PHE:HD2	1:A:254:MET:SD	2.34	0.51
1:A:502:THR:OG1	1:A:505:LYS:HG3	2.10	0.51
1:A:768:ILE:HD11	1:A:1069:TYR:CE1	2.45	0.51
1:A:606:ALA:O	1:A:609:ILE:HG22	2.11	0.51
1:A:344:VAL:CG1	1:A:345:HIS:N	2.73	0.50
1:A:801:LEU:O	1:A:830:VAL:HG13	2.12	0.50
1:A:187:LEU:HD11	1:A:220:GLN:HB2	1.94	0.49
1:A:223:VAL:O	1:A:254:MET:HA	2.13	0.49
1:A:558:LEU:O	1:A:562:LEU:HD12	2.14	0.48
1:A:831:ILE:HD13	1:A:864:LEU:HD13	1.94	0.48
1:A:926:ASN:HA	1:A:1112:PHE:CE2	2.48	0.48
1:A:857:ASP:HA	3:A:1202:B12:H5R2	1.96	0.48
1:A:344:VAL:CG1	1:A:345:HIS:H	2.27	0.47
1:A:353:VAL:HG13	1:A:362:LEU:HD11	1.96	0.47
1:A:531:ASP:HA	1:A:568:VAL:O	2.14	0.47
1:A:749:LEU:HD13	1:A:800:ALA:HB3	1.95	0.47
1:A:919:ALA:HB2	1:A:967:MET:HE1	1.95	0.47
1:A:236:ASP:C	1:A:236:ASP:OD1	2.56	0.47
1:A:322:CYS:O	1:A:323:CYS:C	2.57	0.47
1:A:344:VAL:HG12	1:A:345:HIS:H	1.78	0.47
1:A:237:GLU:HG2	1:A:269:TYR:CD2	2.49	0.47
1:A:499:MET:HB3	1:A:535:PHE:HZ	1.80	0.47
1:A:935:TRP:NE1	1:A:1074:PHE:HA	2.31	0.46
1:A:924:TYR:HA	1:A:1108:ILE:HG21	1.98	0.46
1:A:767:ASP:C	1:A:767:ASP:OD1	2.58	0.46
1:A:1015:LEU:HD23	1:A:1175:PHE:HE2	1.79	0.46
1:A:1016:VAL:HA	1:A:1019:PHE:CE2	2.51	0.46
1:A:535:PHE:HB2	1:A:547:LEU:HD23	1.97	0.46
1:A:758:VAL:HG22	3:A:1202:B12:H421	1.98	0.46
1:A:49:VAL:HG21	1:A:311:VAL:CG2	2.46	0.45
1:A:1094:MET:CE	1:A:1166:VAL:HG21	2.46	0.45
1:A:1151:THR:HG23	1:A:1155:GLN:O	2.17	0.45
1:A:169:LEU:HD21	1:A:198:GLN:HG2	1.98	0.45
1:A:122:LEU:HD12	1:A:122:LEU:HA	1.76	0.45
1:A:533:LEU:HD13	1:A:535:PHE:HZ	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:GLU:HB3	1:A:603:ILE:HG22	1.98	0.45
1:A:804:SER:HB3	3:A:1202:B12:H4B	1.99	0.45
1:A:925:VAL:HG13	1:A:1081:GLU:HB3	1.98	0.45
1:A:332:LYS:HA	1:A:332:LYS:HD2	1.74	0.44
1:A:535:PHE:CB	1:A:547:LEU:HD23	2.47	0.44
1:A:1109:ARG:HA	1:A:1112:PHE:CE2	2.52	0.44
1:A:1131:LEU:HG	1:A:1154:TYR:HB2	2.00	0.44
1:A:762:GLY:O	1:A:766:VAL:HG23	2.17	0.44
1:A:761:ILE:HG12	1:A:1076:HIS:CE1	2.52	0.44
1:A:1028:SER:OG	1:A:1029:ASP:N	2.49	0.44
1:A:464:ASN:HA	1:A:490:VAL:HB	2.00	0.43
1:A:749:LEU:HB2	1:A:865:MET:CE	2.48	0.43
1:A:806:LEU:HD21	3:A:1202:B12:H561	1.99	0.43
1:A:113:LEU:HD11	1:A:173:TYR:CE1	2.53	0.43
1:A:263:MET:HE2	1:A:263:MET:CA	2.48	0.43
1:A:585:VAL:HG21	1:A:644:PHE:HZ	1.83	0.43
1:A:1135:ALA:HB2	1:A:1154:TYR:HE1	1.84	0.43
1:A:531:ASP:O	1:A:569:LEU:HD23	2.19	0.43
1:A:728:MET:SD	1:A:732:VAL:HG23	2.59	0.43
1:A:736:GLU:HB3	1:A:737:PRO:HD3	2.01	0.43
1:A:330:ILE:HA	1:A:333:VAL:HG22	2.00	0.43
1:A:765:LEU:O	1:A:768:ILE:HG22	2.19	0.43
1:A:704:MET:O	1:A:707:VAL:HB	2.18	0.42
1:A:753:THR:HA	1:A:804:SER:HG	1.84	0.42
1:A:759:HIS:HE1	1:A:807:LEU:HD21	1.84	0.42
1:A:1069:TYR:O	1:A:1070:GLN:C	2.61	0.42
1:A:1076:HIS:CG	1:A:1077:GLY:N	2.87	0.42
3:A:1202:B12:H411	3:A:1202:B12:H363	1.53	0.42
1:A:436:VAL:HG12	1:A:438:VAL:H	1.85	0.42
1:A:375:LEU:HG	1:A:402:GLN:HE22	1.85	0.42
1:A:31:ASN:O	1:A:35:ARG:HG2	2.20	0.42
1:A:749:LEU:HD22	1:A:865:MET:SD	2.60	0.42
1:A:766:VAL:HG21	3:A:1202:B12:C7B	2.49	0.42
1:A:902:LEU:HD13	1:A:1043:ASP:OD1	2.18	0.42
1:A:200:ARG:HA	1:A:203:VAL:HG12	2.02	0.42
1:A:375:LEU:HB2	1:A:411:ASP:O	2.20	0.42
1:A:527:ASP:OD1	1:A:527:ASP:C	2.63	0.42
1:A:425:MET:HB3	1:A:425:MET:HE3	1.56	0.42
3:A:1202:B12:H253	3:A:1202:B12:H301	1.82	0.42
1:A:41:TYR:HB2	1:A:186:ASP:OD2	2.20	0.41
1:A:371:VAL:O	1:A:371:VAL:HG12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:GLU:O	1:A:126:ALA:C	2.61	0.41
1:A:1070:GLN:HG2	1:A:1074:PHE:CE2	2.54	0.41
1:A:718:LEU:HB3	1:A:719:PRO:HD3	2.03	0.41
1:A:944:ARG:O	1:A:948:GLU:HG2	2.20	0.41
1:A:1054:GLU:C	1:A:1054:GLU:CD	2.89	0.41
3:A:1202:B12:H91	3:A:1202:B12:H542	1.79	0.41
1:A:263:MET:O	1:A:264:ASP:C	2.63	0.41
1:A:311:VAL:HG13	1:A:312:THR:HG23	2.03	0.41
1:A:953:ARG:HG3	1:A:954:GLU:HG2	2.03	0.41
1:A:976:VAL:HG11	1:A:1090:TRP:CE2	2.55	0.41
1:A:98:LEU:HD13	1:A:144:TYR:CD1	2.56	0.41
1:A:381:LYS:HD2	1:A:381:LYS:H	1.86	0.41
1:A:399:ALA:O	1:A:403:VAL:HG23	2.21	0.41
1:A:584:ARG:HD2	1:A:608:LYS:HB3	2.03	0.41
1:A:963:LEU:O	1:A:967:MET:HG3	2.21	0.41
1:A:130:ALA:HB1	1:A:180:LEU:HD12	2.02	0.40
1:A:392:LEU:HD12	1:A:431:LYS:HB3	2.01	0.40
1:A:692:PRO:HB2	1:A:739:MET:CG	2.51	0.40
1:A:1029:ASP:N	1:A:1029:ASP:OD1	2.54	0.40
1:A:707:VAL:HG22	1:A:716:MET:HE1	2.03	0.40
1:A:1005:PHE:HD2	1:A:1016:VAL:HB	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	1127/1178 (96%)	1093 (97%)	34 (3%)	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	914/949 (96%)	910 (100%)	4 (0%)	84 82

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	LEU
1	A	331	ARG
1	A	857	ASP
1	A	1076	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	773	ASN
1	A	923	HIS
1	A	1048	GLN
1	A	1155	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

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	$\#\|Z\| > 2$
3	B12	A	1202	-	94,101,101	1.00	5 (5%)	149,166,166	1.88	39 (26%)

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
3	B12	A	1202	-	-	24/56/223/223	0/3/11/11

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1202	B12	C14-N23	2.69	1.38	1.35
3	A	1202	B12	C35-C5	2.61	1.56	1.50
3	A	1202	B12	C54-C17	2.52	1.58	1.54
3	A	1202	B12	C2B-N1B	-2.24	1.33	1.37
3	A	1202	B12	C30-C3	2.15	1.59	1.54

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1202	B12	C19-N24-C16	6.20	114.07	107.29
3	A	1202	B12	C16-C15-C14	-5.86	112.33	121.26
3	A	1202	B12	C55-C17-C16	5.49	127.33	116.59
3	A	1202	B12	C15-C16-N24	4.30	128.54	122.42
3	A	1202	B12	C54-C17-C16	-4.20	90.67	112.41
3	A	1202	B12	C13-C12-C11	-4.01	96.49	100.97
3	A	1202	B12	C17-C16-N24	-3.98	105.09	111.17
3	A	1202	B12	C9-N22-C6	3.89	109.96	105.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1202	B12	C20-C1-C19	3.66	112.88	109.35
3	A	1202	B12	C12-C11-C10	3.66	128.11	123.40
3	A	1202	B12	C1P-N59-C57	-3.42	115.35	122.69
3	A	1202	B12	C9-C10-C11	-3.36	121.16	125.97
3	A	1202	B12	C25-C2-C1	-3.21	108.95	113.75
3	A	1202	B12	C5M-C5B-C6B	-3.16	114.31	120.76
3	A	1202	B12	C55-C17-C18	3.11	117.06	111.12
3	A	1202	B12	C25-C2-C26	3.04	115.77	109.74
3	A	1202	B12	C12-C13-C14	-2.98	97.36	102.26
3	A	1202	B12	C12-C11-N23	-2.84	107.91	111.83
3	A	1202	B12	C9B-C8B-N1B	2.81	106.55	105.30
3	A	1202	B12	C7B-C8B-C9B	2.79	125.82	122.47
3	A	1202	B12	C54-C17-C18	-2.76	109.02	112.99
3	A	1202	B12	C41-C8-C7	-2.75	106.62	114.19
3	A	1202	B12	C20-C1-C2	-2.73	108.78	113.28
3	A	1202	B12	C2P-C1P-N59	-2.68	108.99	112.92
3	A	1202	B12	C17-C18-C19	-2.67	98.31	102.36
3	A	1202	B12	C4B-C5B-C6B	2.64	123.56	119.69
3	A	1202	B12	O28-C27-N29	-2.58	115.63	122.53
3	A	1202	B12	C30-C3-C4	2.49	115.50	109.66
3	A	1202	B12	C2-C1-C19	-2.35	114.96	118.61
3	A	1202	B12	C55-C56-C57	-2.20	106.35	111.25
3	A	1202	B12	O2-C3R-C2R	2.18	119.48	111.68
3	A	1202	B12	C2-C3-C4	2.15	104.06	101.64
3	A	1202	B12	C42-C43-N45	2.15	123.37	116.49
3	A	1202	B12	C19-C1-N21	2.15	104.35	102.14
3	A	1202	B12	O6R-C1R-N1B	-2.12	104.01	108.09
3	A	1202	B12	C8B-C9B-N3B	-2.08	107.75	110.00
3	A	1202	B12	C53-C15-C16	2.05	123.84	120.36
3	A	1202	B12	C9B-C4B-C5B	-2.04	117.05	120.83
3	A	1202	B12	C48-C13-C12	-2.03	110.73	116.52

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1202	B12	C1-C2-C26-C27
3	A	1202	B12	C25-C2-C26-C27
3	A	1202	B12	C3-C2-C26-C27
3	A	1202	B12	C12-C13-C48-C49
3	A	1202	B12	C14-C13-C48-C49
3	A	1202	B12	C1P-C2P-O3-P

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Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	A	1202	B12	C3P-C2P-O3-P
3	A	1202	B12	O6R-C4R-C5R-O8R
3	A	1202	B12	C4-C3-C30-C31
3	A	1202	B12	C42-C41-C8-C9
3	A	1202	B12	C3R-C4R-C5R-O8R
3	A	1202	B12	C16-C17-C55-C56
3	A	1202	B12	C38-C37-C7-C6
3	A	1202	B12	C2-C3-C30-C31
3	A	1202	B12	C13-C48-C49-C50
3	A	1202	B12	C2R-C1R-N1B-C8B
3	A	1202	B12	O6R-C1R-N1B-C8B
3	A	1202	B12	C17-C18-C60-C61
3	A	1202	B12	C38-C37-C7-C8
3	A	1202	B12	C2R-C1R-N1B-C2B
3	A	1202	B12	C19-C18-C60-C61
3	A	1202	B12	C2P-O3-P-O4
3	A	1202	B12	C2P-O3-P-O5
3	A	1202	B12	C2P-C1P-N59-C57

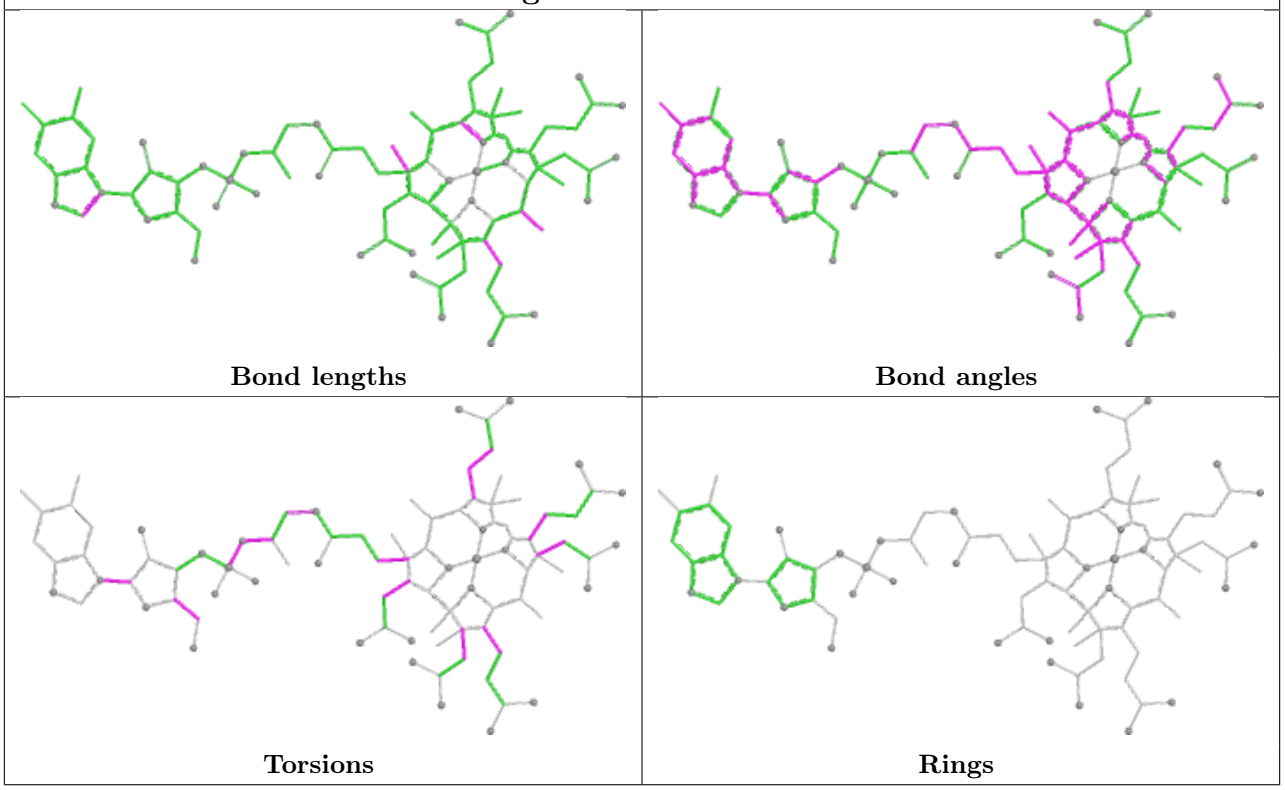
There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1202	B12	13	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 then 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.

Ligand B12 A 1202



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.

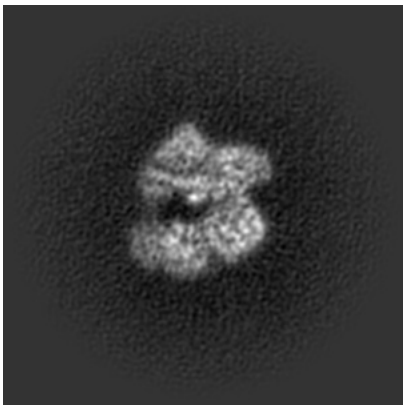
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-78314. These allow visual inspection of the internal detail of the map and identification of artifacts.

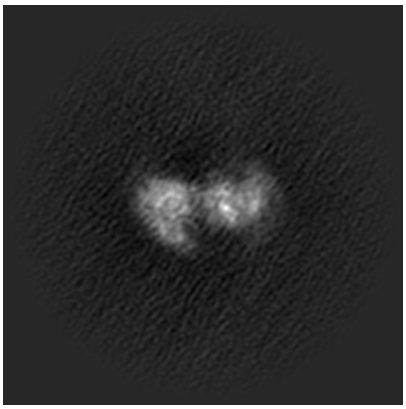
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

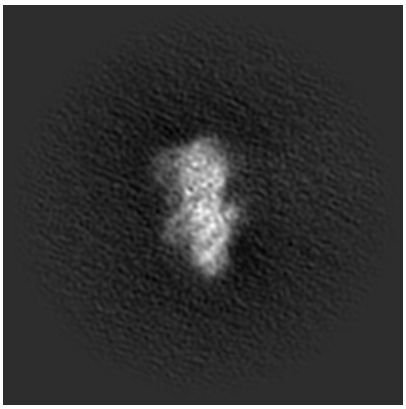
6.1.1 Primary map



X

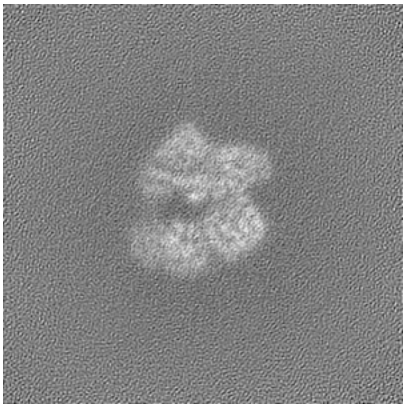


Y

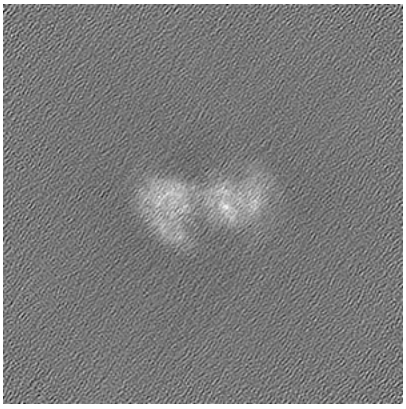


Z

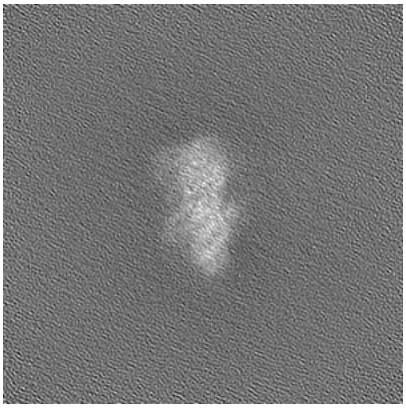
6.1.2 Raw map



X



Y

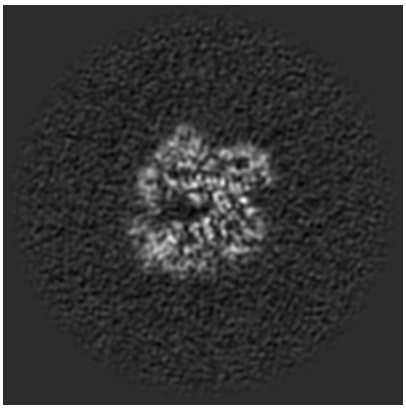


Z

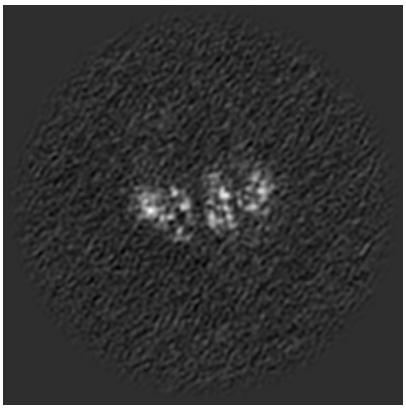
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

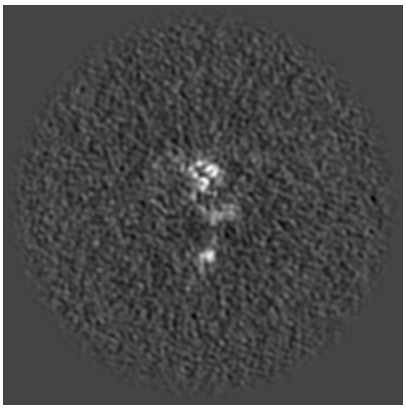
6.2.1 Primary map



X Index: 128

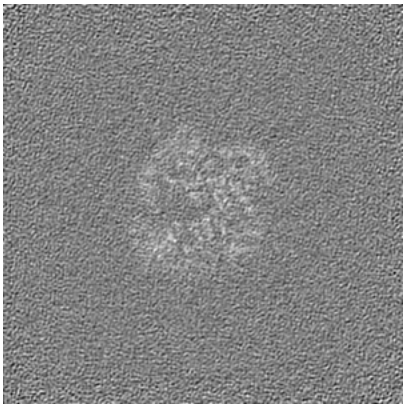


Y Index: 128

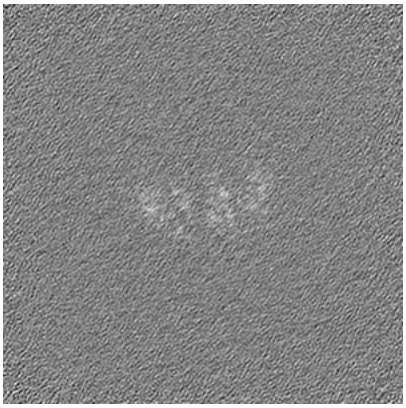


Z Index: 128

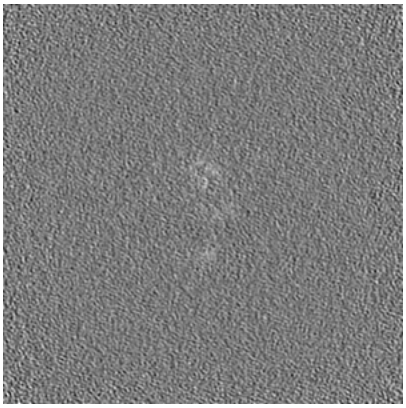
6.2.2 Raw map



X Index: 128



Y Index: 128

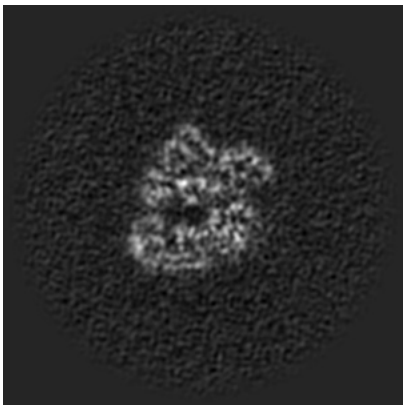


Z Index: 128

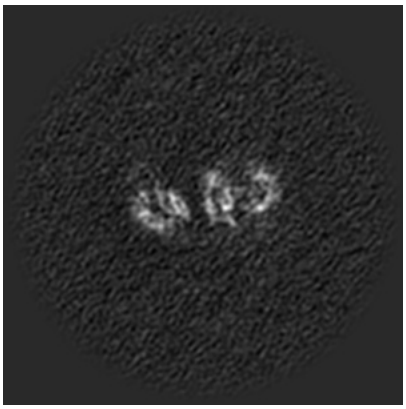
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

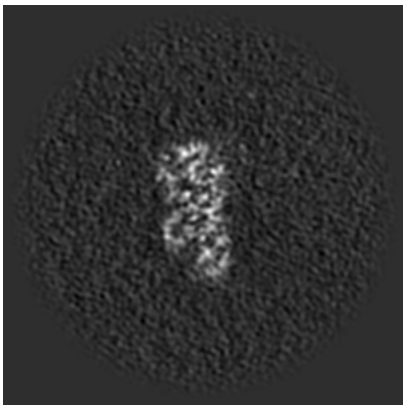
6.3.1 Primary map



X Index: 134

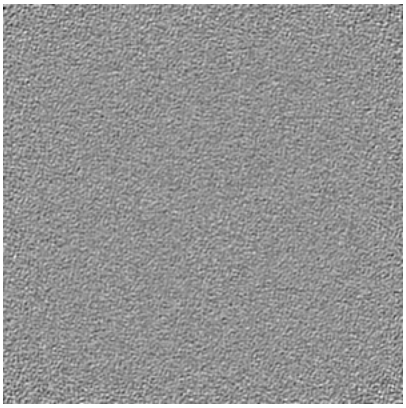


Y Index: 124

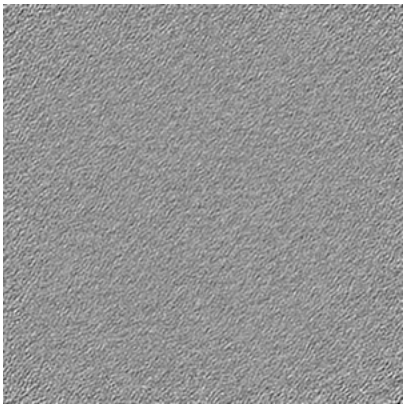


Z Index: 108

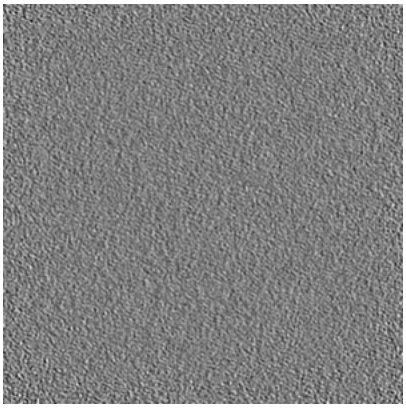
6.3.2 Raw map



X Index: 0



Y Index: 0

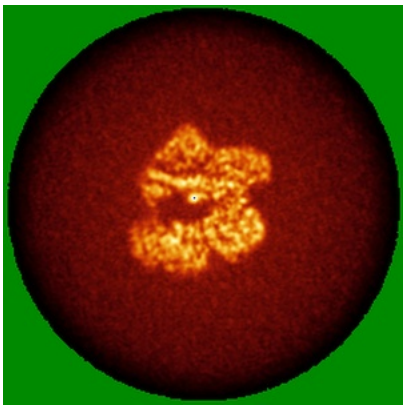


Z Index: 0

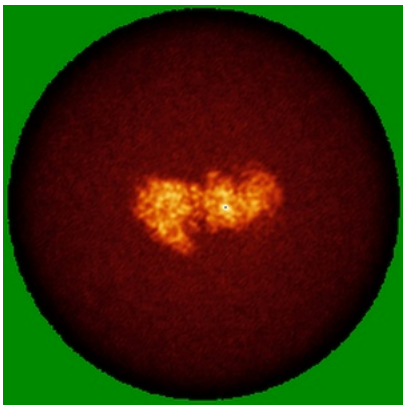
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

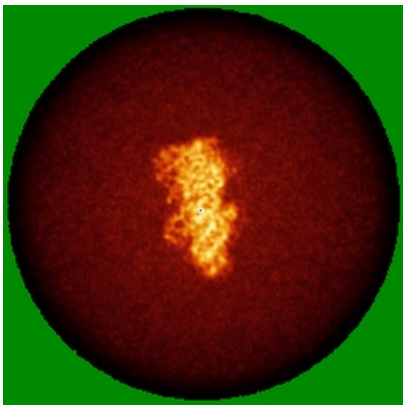
6.4.1 Primary map



X

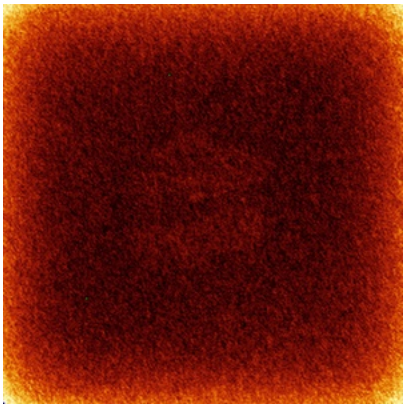


Y

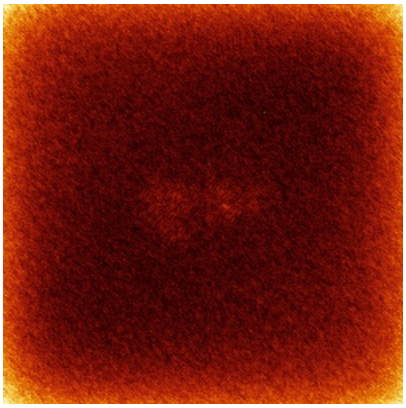


Z

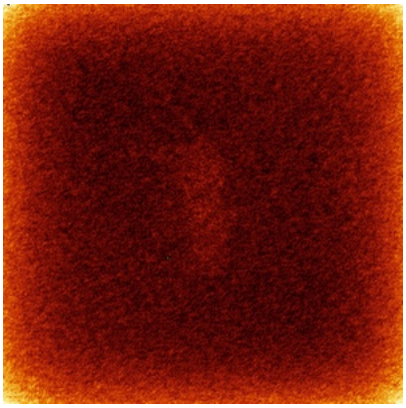
6.4.2 Raw map



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 0.119. 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.

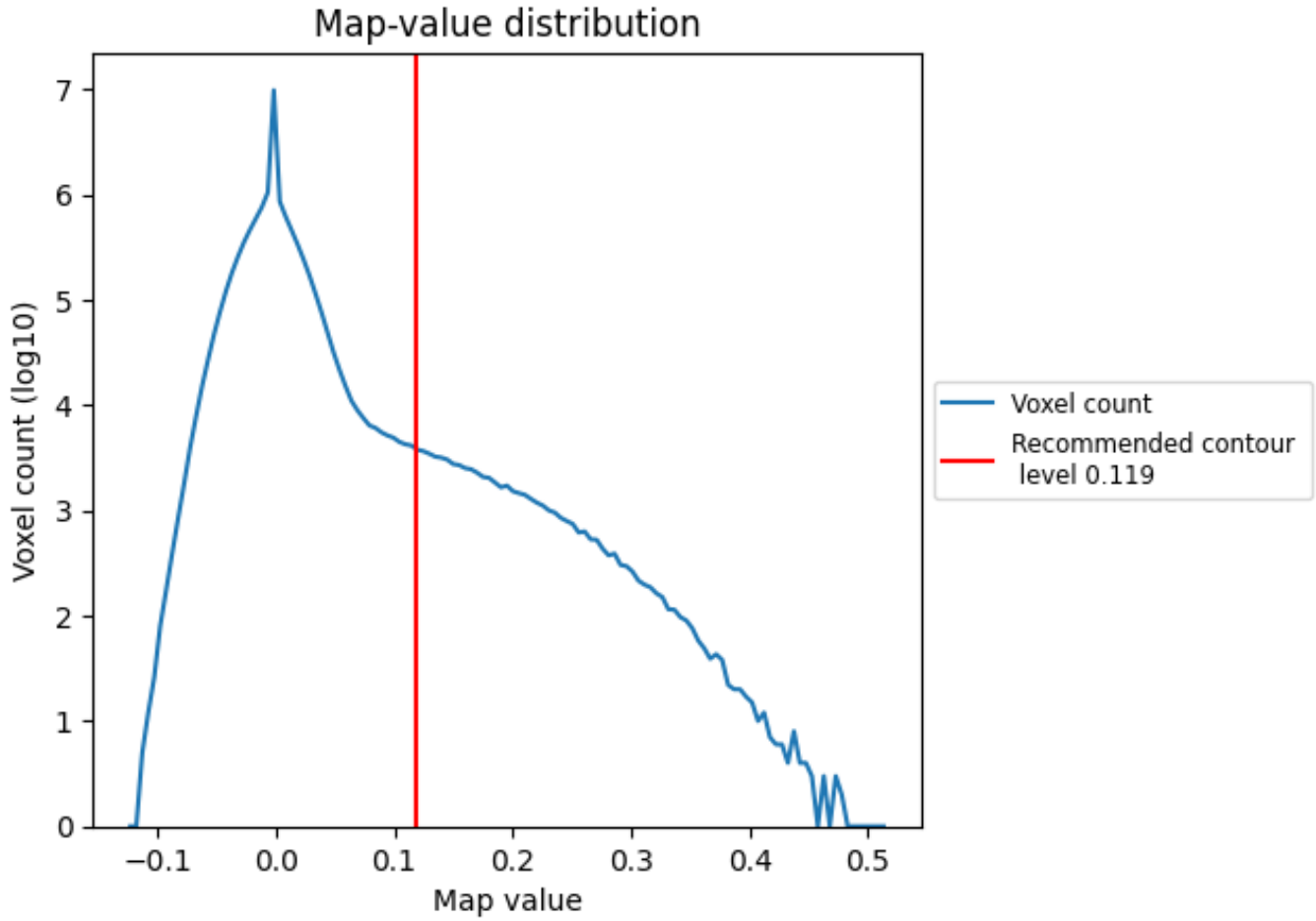
6.6 Mask visualisation (i)

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

7 Map analysis

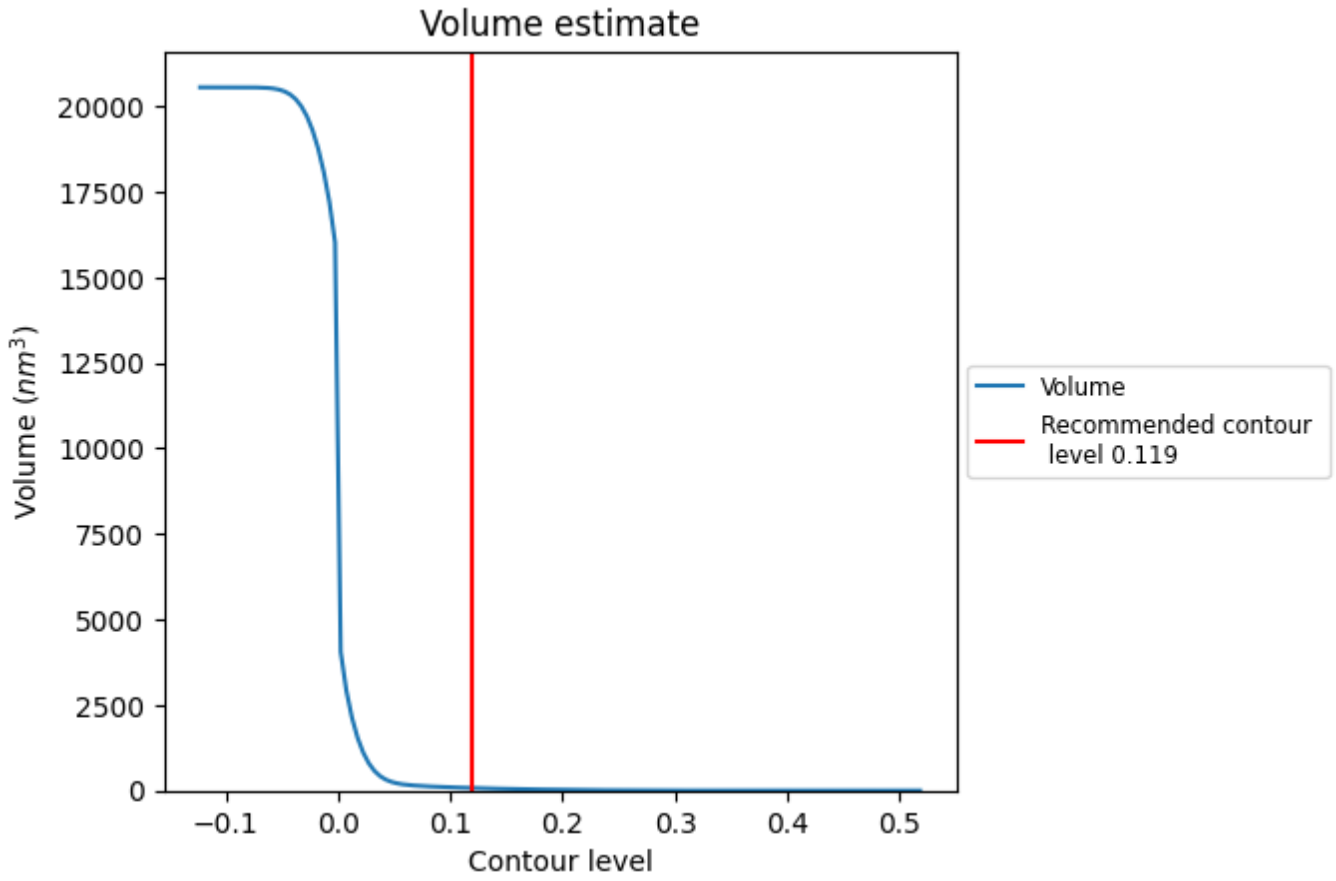
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

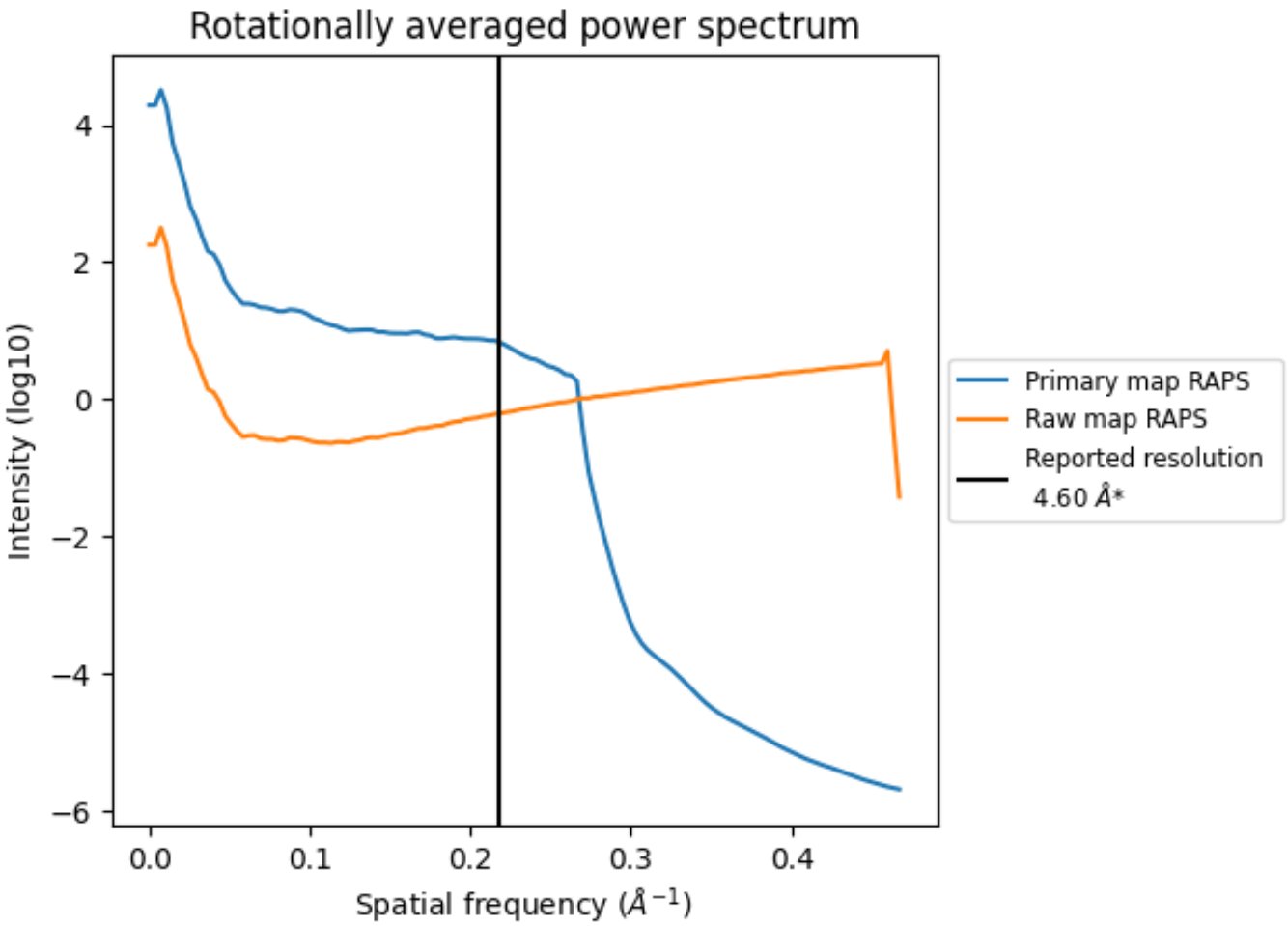
7.2 Volume estimate



The volume at the recommended contour level is 75 nm^3; this corresponds to an approximate mass of 68 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum (i)

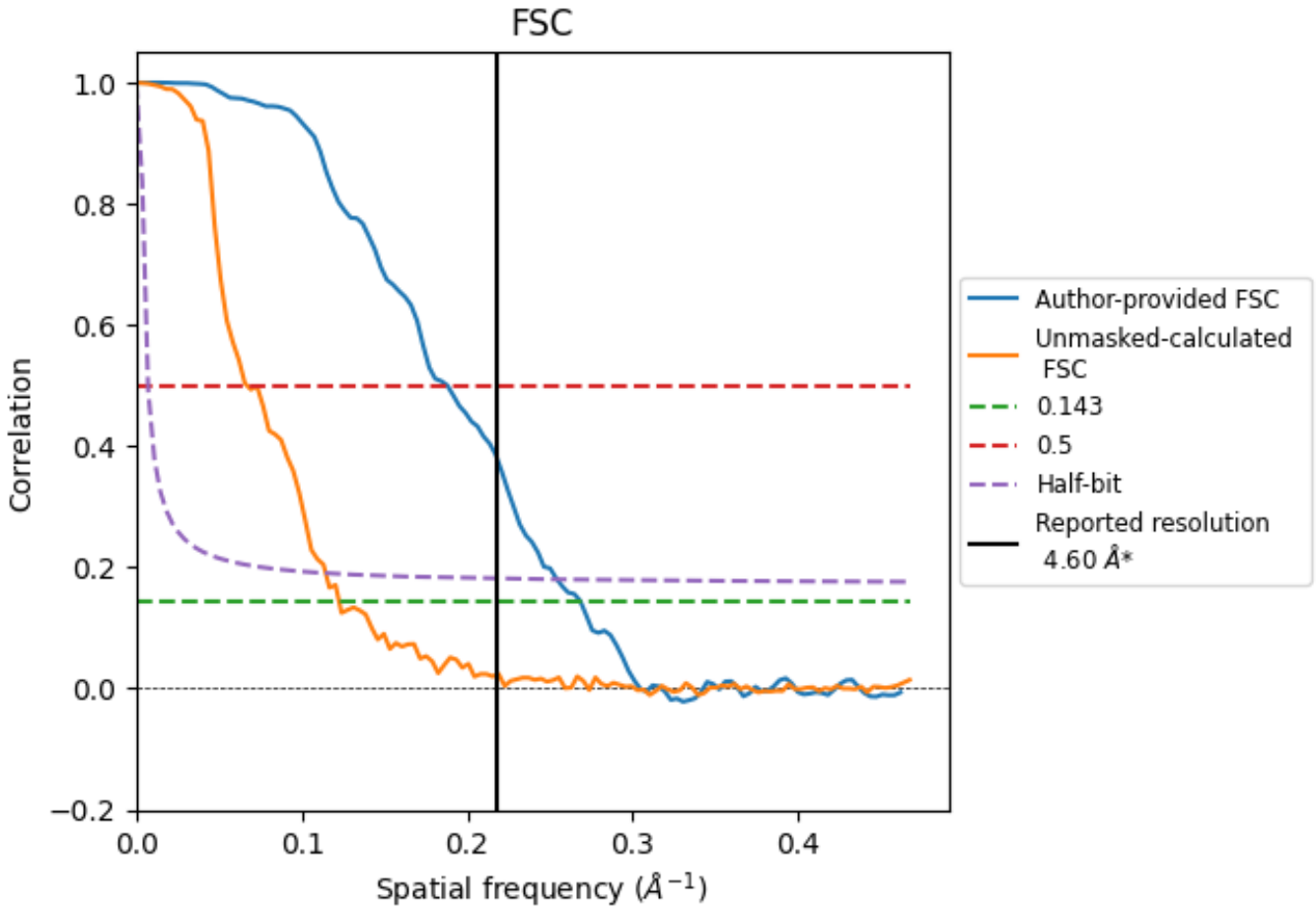


*Reported resolution corresponds to spatial frequency of 0.217 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.217 Å^{-1}

8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.60	-	-
Author-provided FSC curve	3.72	5.32	3.93
Unmasked-calculated*	8.15	14.93	8.73

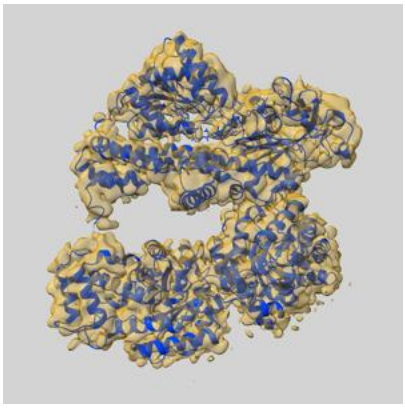
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 3.72 differs from the reported value 4.6 by more than 10 %

The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 8.15 differs from the reported value 4.6 by more than 10 %

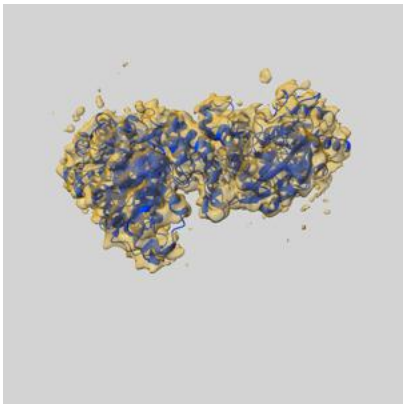
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-78314 and PDB model 37MQ. Per-residue inclusion information can be found in section 3 on page 4.

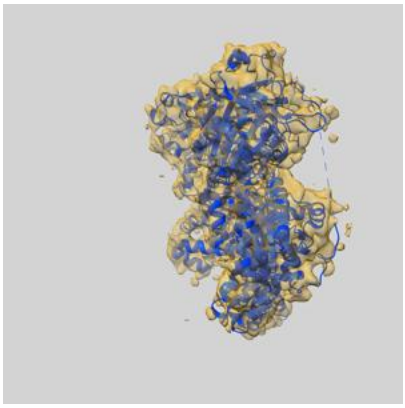
9.1 Map-model overlay



X



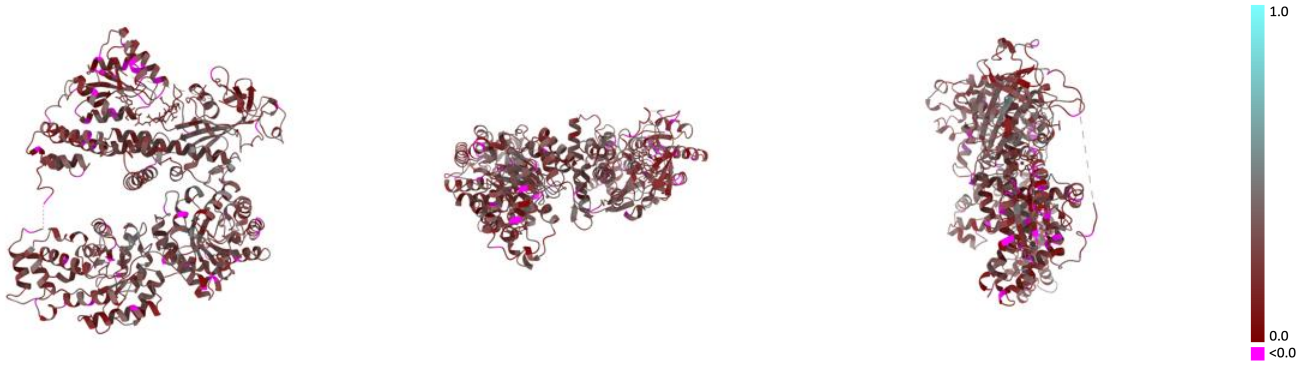
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.119 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.

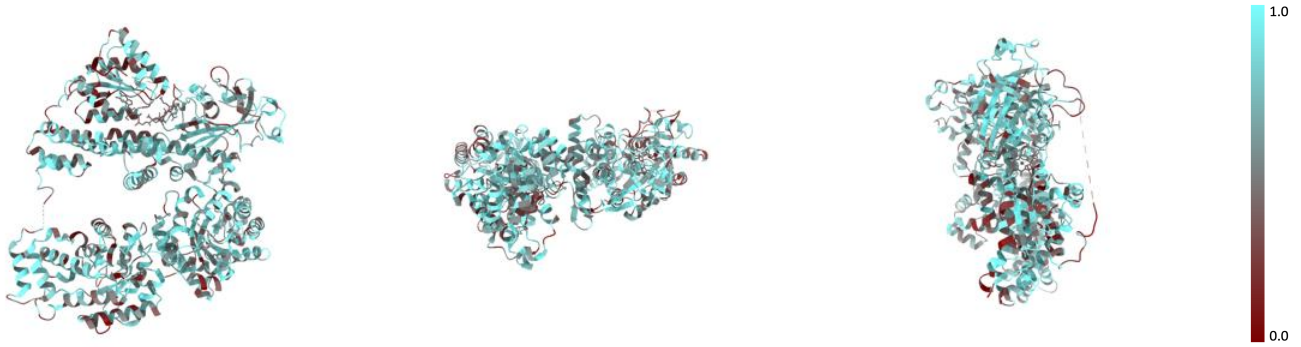
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

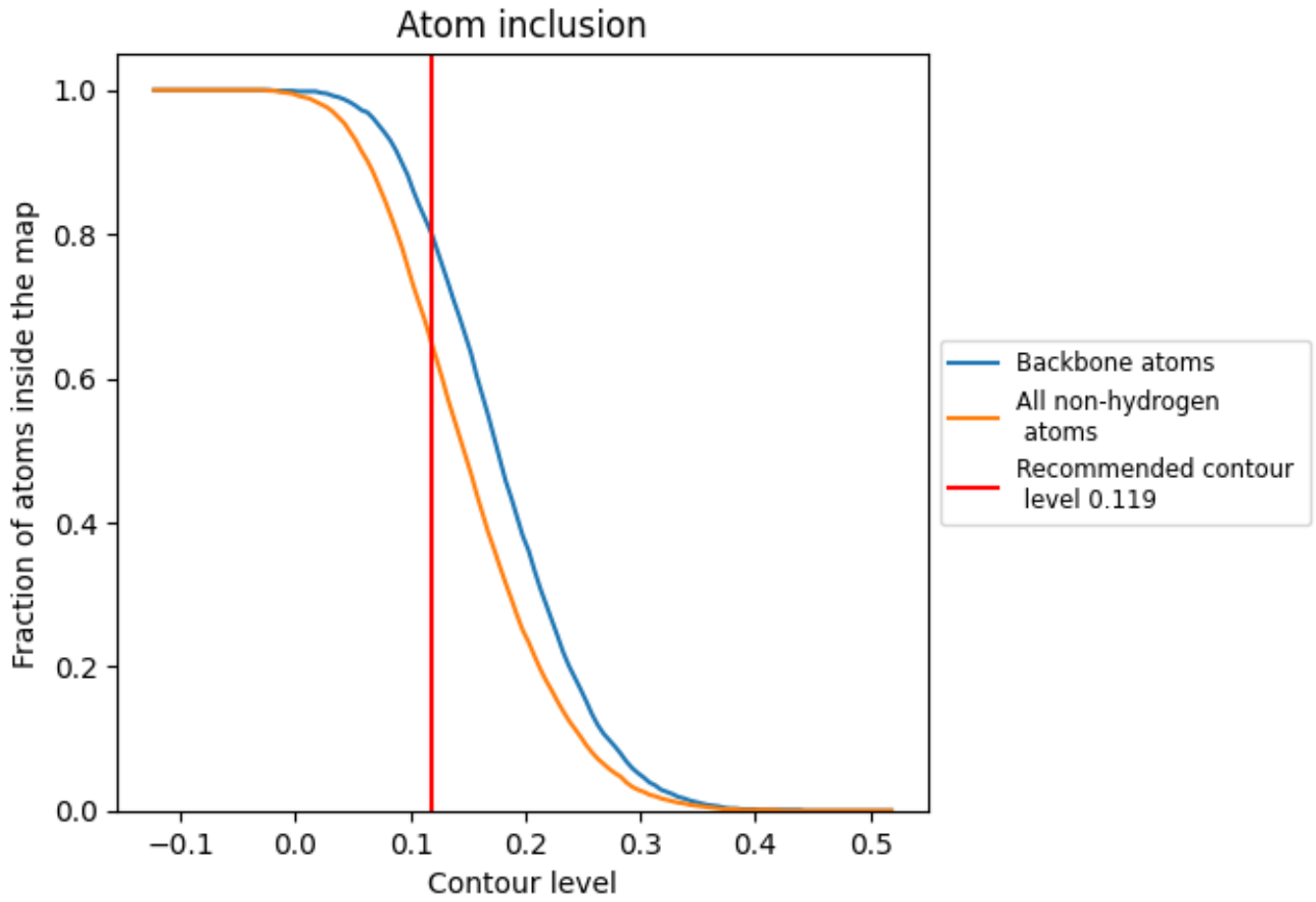
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.119).

9.4 Atom inclusion (i)



At the recommended contour level, 80% of all backbone atoms, 64% 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 (0.119) and Q-score for the entire model and for each chain.



Chain	Atom inclusion	Q-score
All	0.6450	0.2510
A	0.6450	0.2510