Full wwPDB EM Validation Report (i)

Aug 26, 2026 – 01:53 pm BST

PDB ID	:	9SSS / pdb_00009sss
EMDB ID	:	EMD-55193
Title	:	Human Methionine Synthase With Methyltetrahydrofolate, Hydroxocobalamin, and SAM, C-Half His-ON From Full-Length
Authors	:	Ferreira, D.S.M.; Yue, W.W.; McCorvie, T.J.
Deposited on	:	2025-09-26
Resolution	:	3.01 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 (i)) were used in the production of this report:

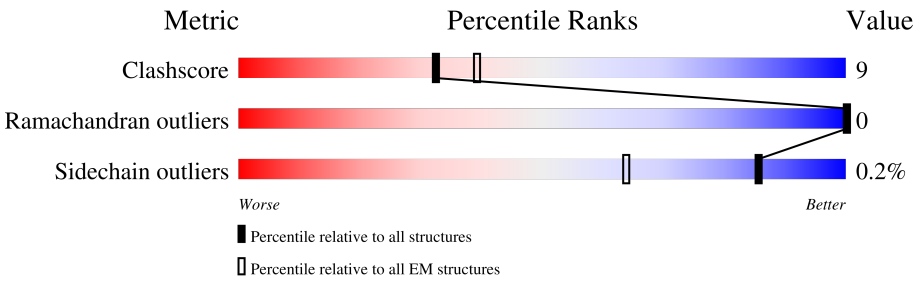
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	1.8.4, CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data 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)	:	2.50

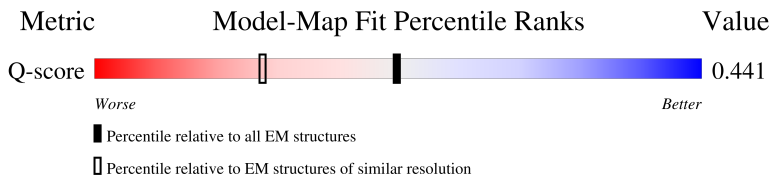
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.01 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	13882 (2.51 - 3.51)

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	1265	9% 38% 10% 52%

2 Entry composition (i)

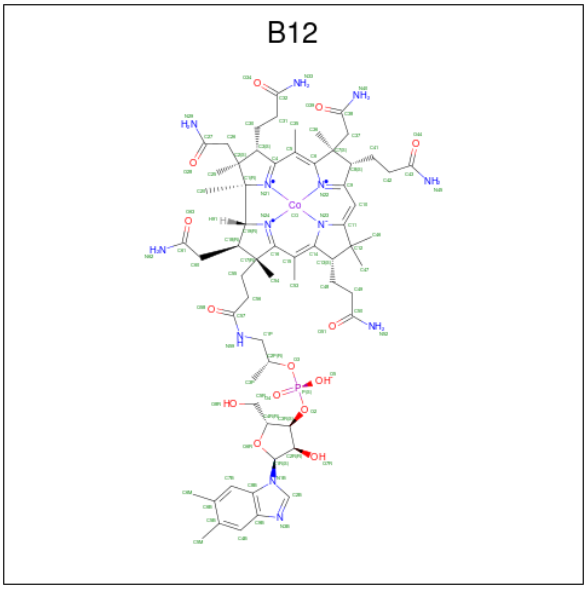
There are 2 unique types of molecules in this entry. The entry contains 4876 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	601	4785	3054	811	900	20	0	0

- Molecule 2 is COBALAMIN (CCD ID: B12) (formula: $C_{62}H_{89}CoN_{13}O_{14}P$) (labeled as "Ligand of Interest" by depositor).



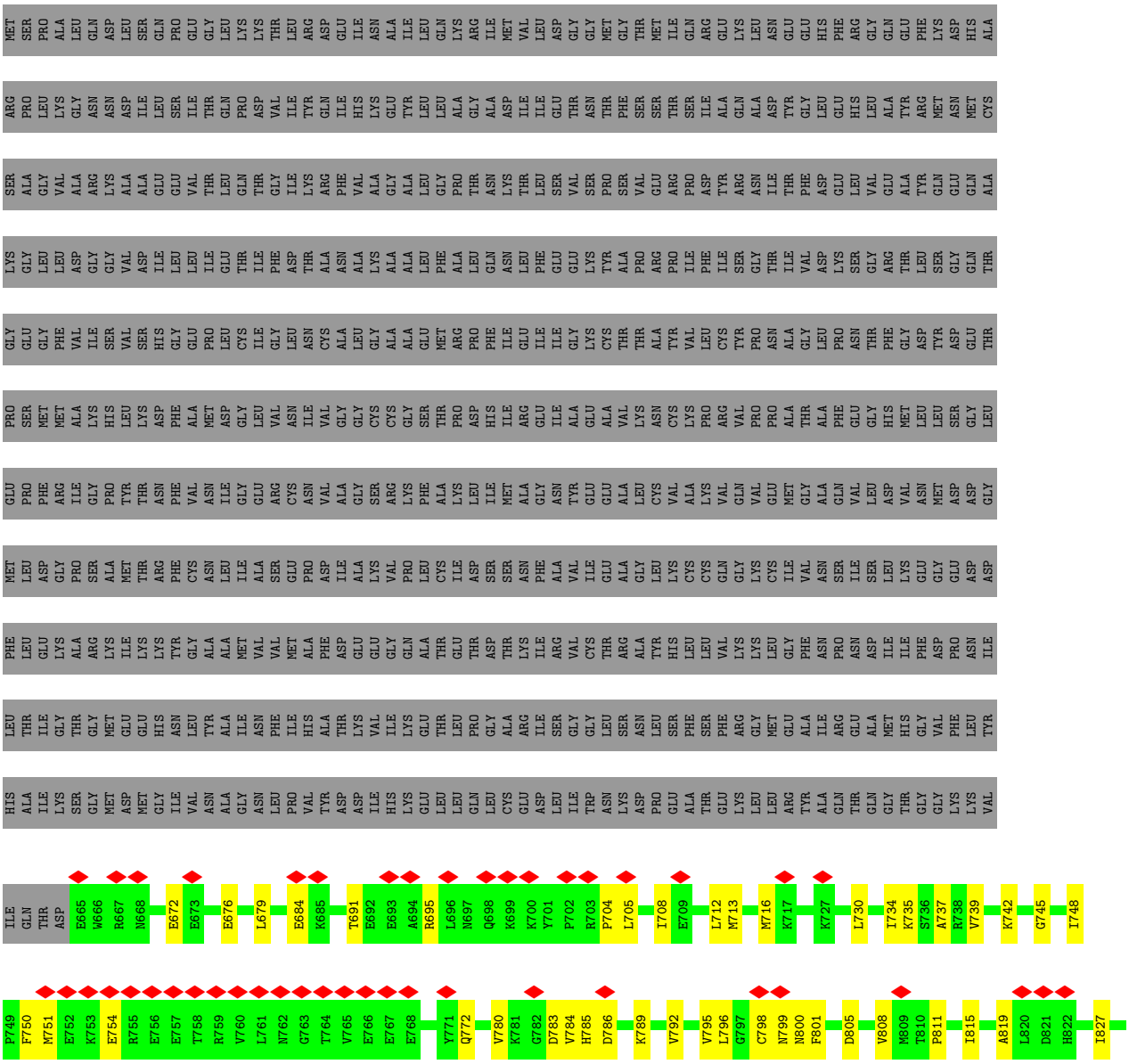
Mol	Chain	Residues	Atoms						AltConf
			Total	C	Co	N	O	P	
2	A	1	91	62	1	13	14	1	0

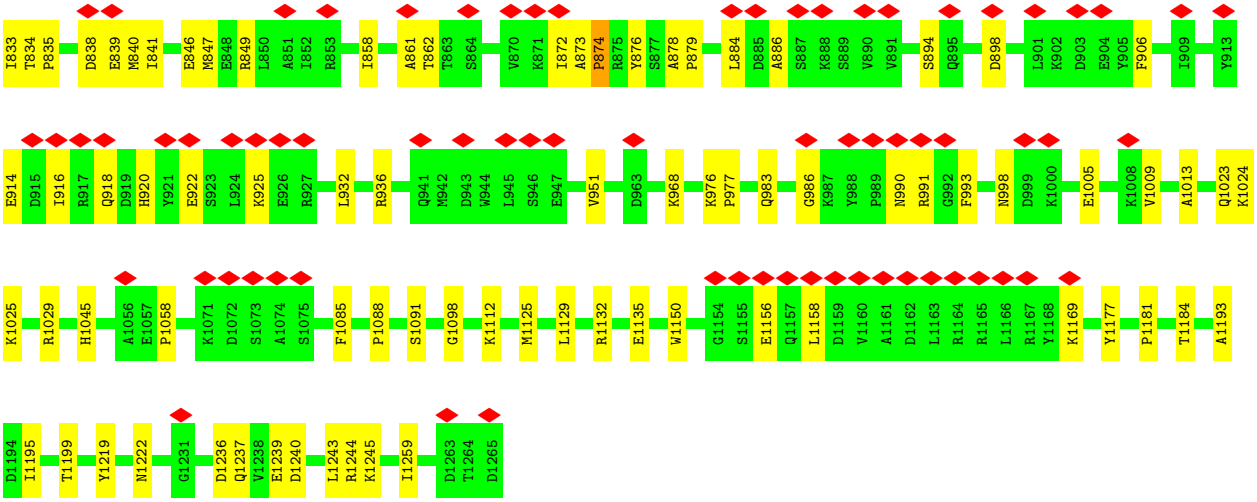
3 Residue-property plots (i)

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:




I833 T834 P835 D838 E839 M840 I841 E846 M847 E848 R849 L850 A851 I852 R853 I858 A861 T862 T863 S864 V870 K871 I872 A873 P874 R875 Y876 S877 A878 P879 L884 D885 A886 S887 K888 S889 V890 V891 S894 Q895 D898 L901 K902 D903 E904 Y905 F906 I909 Y913
E914 D915 I916 R917 Q918 D919 H920 Y921 E922 S923 L924 K925 E926 R927 L932 R936 Q941 M942 D943 W944 L945 S946 E947 V951 D963 K968 K976 P977 Q983 G986 K987 Y988 P989 N990 R991 G992 F993 N998 D999 K1000 E1005 K1008 V1009 A1013 Q1023 K1024
K1025 R1029 H1045 A1056 E1057 P1058 K1071 D1072 S1073 A1074 S1075 F1085 P1088 S1091 G1098 K1112 M1125 L1129 R1132 E1135 W1150 G1154 S1155 E1156 Q1157 L1158 D1159 V1160 A1161 D1162 L1163 R1164 R1165 L1166 R1167 Y1168 K1169 Y1177 P1181 T1184 A1193
D1194 I1195 T1199 Y1219 N1222 G1231 D1236 Q1237 Y1238 E1239 D1240 L1243 R1244 K1245 I1259 D1263 T1264 D1265

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	25594	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 ($e^-/Å^2$)	65.2	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	165000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.470	Depositor
Minimum map value	-0.241	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.08	Depositor
Map size (Å)	203.20001, 203.20001, 203.20001	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.016, 1.016, 1.016	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: 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	0.13	0/4890	0.37	2/6615 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	874	PRO	CA-N-CD	-6.22	103.29	112.00
1	A	862	THR	N-CA-C	-5.11	108.32	114.75

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	4785	0	4769	78	0
2	A	91	0	88	13	0
All	All	4876	0	4857	83	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 9.

All (83) 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:745:GLY:HA2	1:A:748:ILE:HD12	1.66	0.77
1:A:684:GLU:HG3	1:A:739:VAL:HG22	1.73	0.70
1:A:833:ILE:HG22	2:A:1301:B12:H302	1.74	0.70
1:A:1184:THR:OG1	1:A:1245:LYS:NZ	2.27	0.67
2:A:1301:B12:H552	2:A:1301:B12:H531	1.81	0.63
1:A:983:GLN:NE2	2:A:1301:B12:O5	2.33	0.61
1:A:1013:ALA:HB1	1:A:1129:LEU:HD21	1.82	0.61
1:A:786:ASP:H	2:A:1301:B12:H521	1.48	0.60
1:A:796:LEU:HB3	1:A:801:PHE:HB2	1.83	0.58
1:A:713:MET:SD	1:A:800:ASN:ND2	2.76	0.58
1:A:783:ASP:OD1	1:A:785:HIS:ND1	2.30	0.58
1:A:691:THR:O	1:A:695:ARG:HG2	2.03	0.57
1:A:708:ILE:HA	1:A:712:LEU:HD12	1.85	0.57
1:A:916:ILE:O	1:A:920:HIS:ND1	2.37	0.57
1:A:841:ILE:HG22	1:A:872:ILE:HG12	1.86	0.56
1:A:990:ASN:ND2	1:A:998:ASN:OD1	2.39	0.55
1:A:780:VAL:HB	1:A:839:GLU:HG2	1.88	0.55
1:A:886:ALA:H	2:A:1301:B12:H5R2	1.72	0.54
1:A:873:ALA:HB3	1:A:874:PRO:HD2	1.90	0.54
1:A:704:PRO:HD2	1:A:754:GLU:HG3	1.90	0.54
1:A:884:LEU:H	1:A:884:LEU:HD23	1.74	0.52
1:A:986:GLY:O	1:A:991:ARG:NH1	2.43	0.51
1:A:737:ALA:HB2	1:A:798:CYS:SG	2.50	0.51
1:A:1088:PRO:O	1:A:1091:SER:OG	2.28	0.51
1:A:936:ARG:NH2	1:A:1240:ASP:OD1	2.45	0.50
1:A:932:LEU:HD22	1:A:1237:GLN:HA	1.93	0.50
1:A:894:SER:O	1:A:898:ASP:N	2.39	0.50
1:A:861:ALA:N	2:A:1301:B12:O7R	2.44	0.50
1:A:1005:GLU:O	1:A:1009:VAL:HG22	2.11	0.50
1:A:1193:ALA:HB3	1:A:1195:ILE:HG13	1.94	0.49
1:A:1098:GLY:O	1:A:1219:TYR:N	2.42	0.49
1:A:789:LYS:NZ	1:A:805:ASP:OD2	2.31	0.48
1:A:922:GLU:O	1:A:925:LYS:HG2	2.14	0.48
1:A:1177:TYR:HE1	2:A:1301:B12:H372	1.78	0.47
1:A:672:GLU:O	1:A:676:GLU:HG2	2.14	0.47
1:A:750:PHE:HB2	1:A:751:MET:HE2	1.98	0.46
1:A:1045:HIS:HB3	1:A:1058:PRO:HB3	1.98	0.46
1:A:876:TYR:CZ	1:A:878:ALA:HB3	2.51	0.45
1:A:1181:PRO:HG3	1:A:1259:ILE:HG21	1.97	0.45
2:A:1301:B12:H601	2:A:1301:B12:H262	1.98	0.45
2:A:1301:B12:HM53	2:A:1301:B12:HM61	1.74	0.45
1:A:834:THR:HB	1:A:835:PRO:HD3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:784:VAL:HG12	1:A:808:VAL:HG11	2.00	0.44
1:A:838:ASP:HA	1:A:841:ILE:HG12	2.00	0.44
1:A:1024:LYS:HE3	1:A:1024:LYS:HB3	1.86	0.44
1:A:811:PRO:O	1:A:815:ILE:HG13	2.18	0.43
1:A:730:LEU:O	1:A:734:ILE:HG12	2.17	0.43
1:A:796:LEU:HD22	1:A:801:PHE:HD2	1.82	0.43
1:A:1132:ARG:NH1	1:A:1135:GLU:OE1	2.48	0.43
1:A:1156:GLU:HG2	1:A:1158:LEU:H	1.83	0.43
2:A:1301:B12:H8	2:A:1301:B12:H401	1.84	0.43
1:A:819:ALA:HB1	1:A:827:ILE:HD11	2.01	0.43
1:A:879:PRO:HB3	1:A:906:PHE:HB2	2.01	0.43
1:A:846:GLU:CD	1:A:849:ARG:HH12	2.26	0.43
1:A:914:GLU:O	1:A:918:GLN:HG2	2.18	0.43
1:A:742:LYS:HA	1:A:742:LYS:HD2	1.87	0.43
1:A:847:MET:HG2	1:A:876:TYR:CZ	2.54	0.43
1:A:1125:MET:HE2	1:A:1125:MET:HB2	1.88	0.43
1:A:977:PRO:HB3	2:A:1301:B12:H541	2.01	0.42
1:A:716:MET:HG3	1:A:798:CYS:O	2.20	0.42
1:A:796:LEU:O	1:A:801:PHE:N	2.49	0.42
1:A:1023:GLN:HB3	1:A:1025:LYS:HG2	2.00	0.42
1:A:1243:LEU:HD12	1:A:1244:ARG:N	2.34	0.42
1:A:1112:LYS:HE2	1:A:1112:LYS:HA	2.01	0.42
1:A:840:MET:HE1	1:A:858:ILE:HG23	2.02	0.42
2:A:1301:B12:H262	2:A:1301:B12:H91	1.79	0.42
1:A:705:LEU:HD12	1:A:708:ILE:HD11	2.01	0.41
1:A:968:LYS:HB2	1:A:968:LYS:HE2	1.87	0.41
1:A:792:VAL:HG21	2:A:1301:B12:C7B	2.50	0.41
1:A:1029:ARG:NH1	1:A:1199:THR:O	2.52	0.41
1:A:1169:LYS:HB2	1:A:1169:LYS:HE2	1.77	0.41
1:A:876:TYR:CE2	1:A:878:ALA:HB3	2.56	0.41
1:A:951:VAL:O	1:A:1150:TRP:NE1	2.49	0.41
1:A:783:ASP:CG	1:A:785:HIS:HD1	2.24	0.41
1:A:735:LYS:O	1:A:739:VAL:HG23	2.20	0.41
1:A:858:ILE:HD13	1:A:858:ILE:HA	1.85	0.41
1:A:679:LEU:HD11	1:A:712:LEU:HD23	2.02	0.41
1:A:1150:TRP:HD1	1:A:1150:TRP:O	2.03	0.41
1:A:879:PRO:HG3	1:A:906:PHE:CG	2.55	0.40
1:A:795:VAL:O	1:A:799:ASN:N	2.52	0.40
1:A:1085:PHE:HA	1:A:1222:ASN:HD22	1.86	0.40
1:A:976:LYS:HG2	1:A:993:PHE:HB3	2.02	0.40
1:A:1236:ASP:O	1:A:1239:GLU:HG2	2.20	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	599/1265 (47%)	586 (98%)	13 (2%)	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	512/1059 (48%)	511 (100%)	1 (0%)	87 91

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	772	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

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	B12	A	1301	-	93,101,101	0.99	5 (5%)	146,166,166	1.59	24 (16%)

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	B12	A	1301	-	-	20/56/223/223	0/3/11/11

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	B12	C35-C5	2.61	1.56	1.50
2	A	1301	B12	C54-C17	2.57	1.59	1.54
2	A	1301	B12	C14-N23	2.48	1.38	1.35
2	A	1301	B12	C2B-N1B	-2.05	1.33	1.37
2	A	1301	B12	C30-C3	2.05	1.59	1.54

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1301	B12	C16-C15-C14	-4.65	114.19	121.25
2	A	1301	B12	C55-C17-C16	4.58	125.69	116.65
2	A	1301	B12	C54-C17-C18	-4.47	106.38	112.98
2	A	1301	B12	C18-C17-C16	4.40	106.02	100.67
2	A	1301	B12	C54-C17-C16	-3.80	92.67	112.40
2	A	1301	B12	C17-C16-N24	-3.43	105.87	111.15
2	A	1301	B12	C53-C15-C16	3.35	126.14	120.38
2	A	1301	B12	C15-C16-N24	3.27	127.13	122.42
2	A	1301	B12	C1P-N59-C57	-3.22	115.67	122.69
2	A	1301	B12	C5M-C5B-C6B	-2.89	114.82	120.74
2	A	1301	B12	C2P-C1P-N59	-2.87	108.70	112.93
2	A	1301	B12	C9B-C8B-N1B	2.87	106.56	105.35
2	A	1301	B12	C4B-C5B-C6B	2.85	123.75	119.67
2	A	1301	B12	C7B-C8B-C9B	2.78	125.81	122.52
2	A	1301	B12	C55-C56-C57	-2.74	105.25	111.23
2	A	1301	B12	C18-C19-N24	2.67	106.37	102.31
2	A	1301	B12	C37-C7-C8	-2.61	101.39	108.39
2	A	1301	B12	C13-C14-C15	-2.55	120.42	124.32
2	A	1301	B12	C1-C19-N24	-2.52	103.40	106.24
2	A	1301	B12	O28-C27-N29	-2.33	116.16	122.50
2	A	1301	B12	C9B-C4B-C5B	-2.18	116.70	120.71
2	A	1301	B12	C8B-C9B-N3B	-2.11	107.64	109.97
2	A	1301	B12	O2-C3R-C2R	2.06	119.15	111.68
2	A	1301	B12	C31-C32-N33	2.04	122.86	116.51

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1301	B12	C16-C17-C55-C56
2	A	1301	B12	N59-C1P-C2P-C3P
2	A	1301	B12	N59-C1P-C2P-O3
2	A	1301	B12	C1P-C2P-O3-P
2	A	1301	B12	C3P-C2P-O3-P
2	A	1301	B12	C2R-C1R-N1B-C2B
2	A	1301	B12	O6R-C4R-C5R-O8R
2	A	1301	B12	C3R-C4R-C5R-O8R
2	A	1301	B12	C3-C30-C31-C32
2	A	1301	B12	C12-C13-C48-C49
2	A	1301	B12	C18-C17-C55-C56
2	A	1301	B12	C14-C13-C48-C49
2	A	1301	B12	C2R-C1R-N1B-C8B
2	A	1301	B12	C17-C55-C56-C57

Continued on next page...

Continued from previous page...

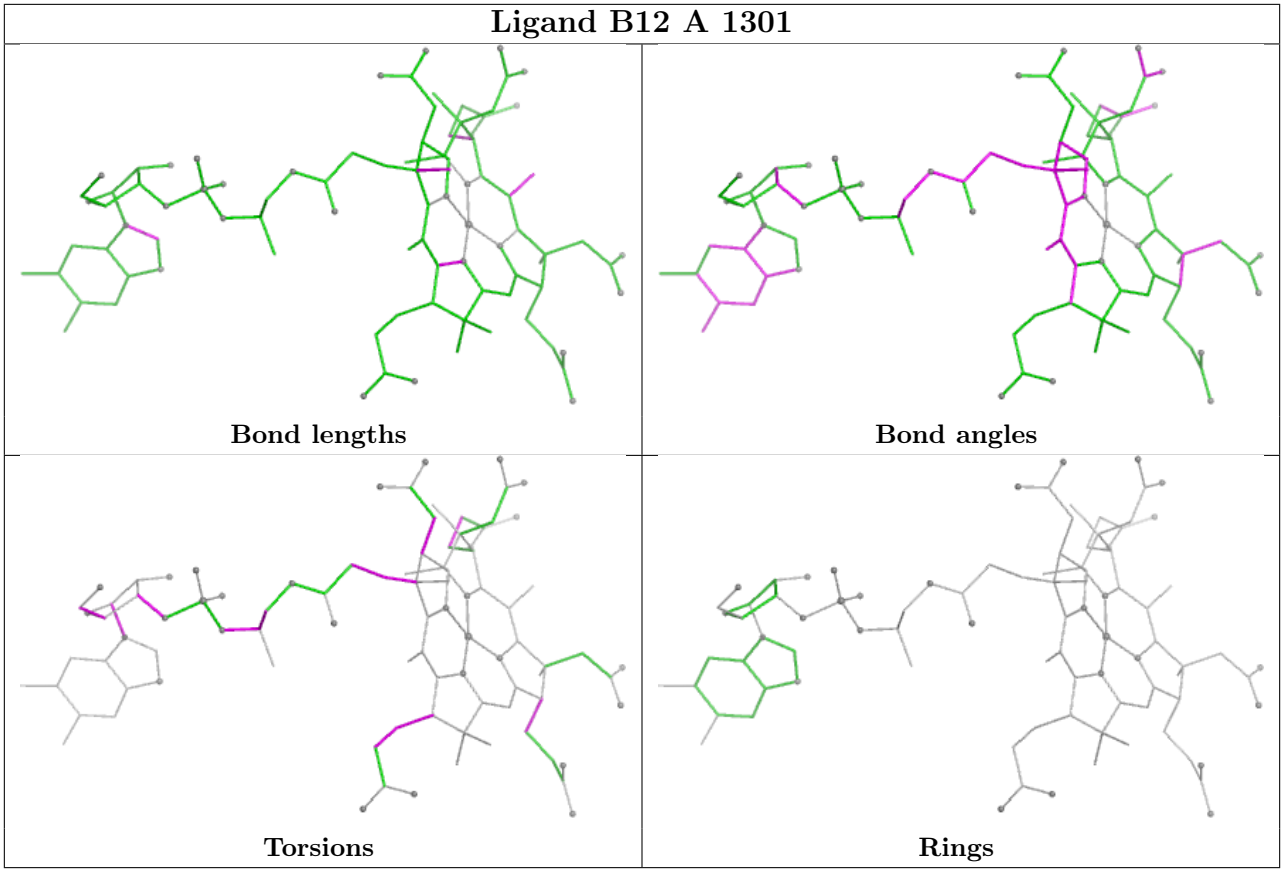
Mol	Chain	Res	Type	Atoms
2	A	1301	B12	C2R-C3R-O2-P
2	A	1301	B12	O6R-C1R-N1B-C8B
2	A	1301	B12	C42-C41-C8-C7
2	A	1301	B12	C19-C18-C60-C61
2	A	1301	B12	C13-C48-C49-C50
2	A	1301	B12	C4R-C3R-O2-P

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1301	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.



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.

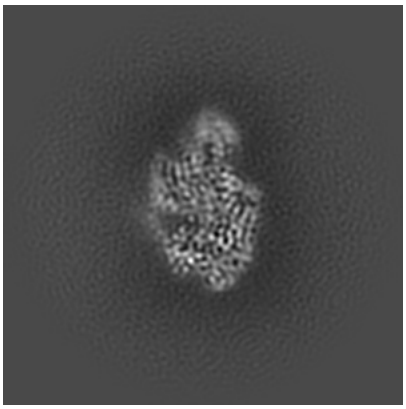
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-55193. These allow visual inspection of the internal detail of the map and identification of artifacts.

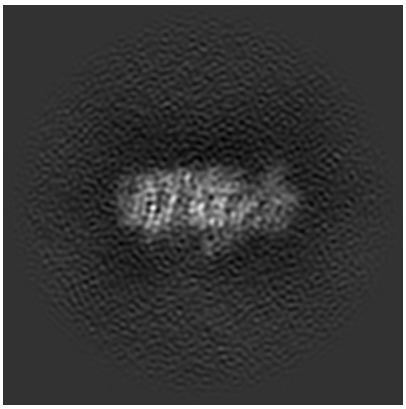
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

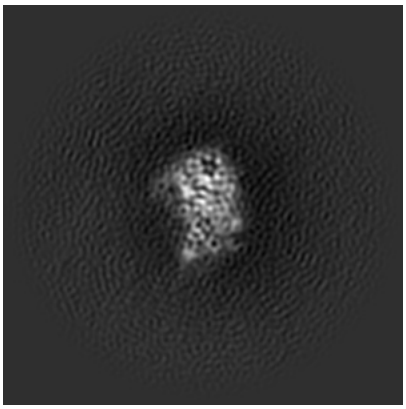
6.1.1 Primary map



X

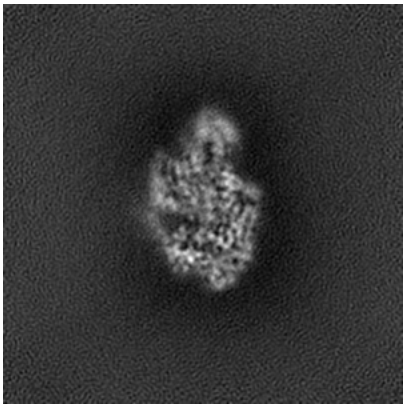


Y

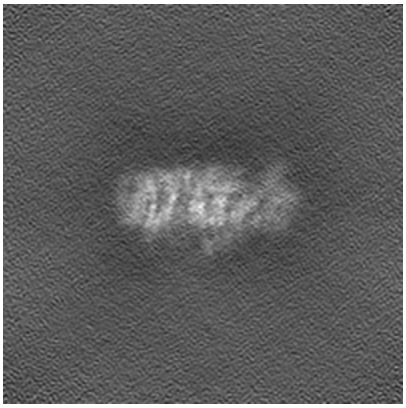


Z

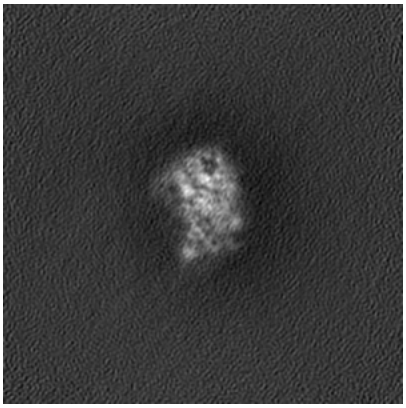
6.1.2 Raw map



X



Y

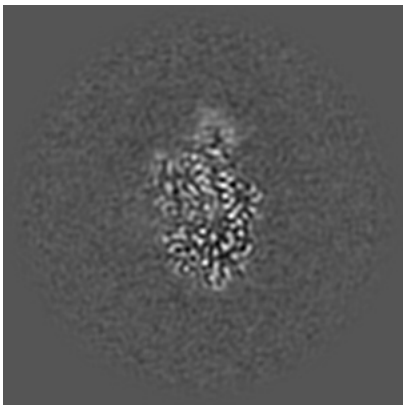


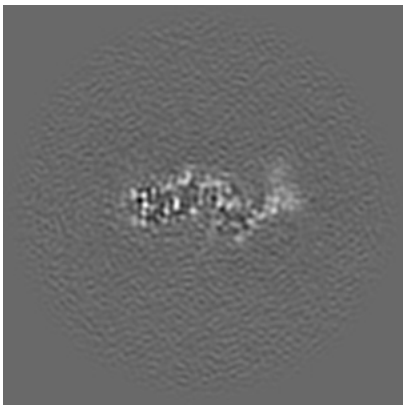
Z

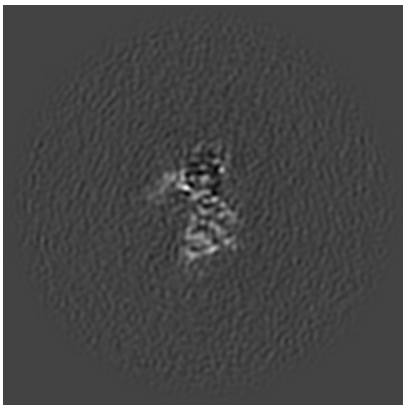
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

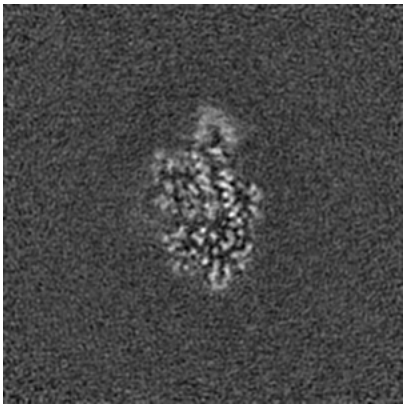
6.2.1 Primary map

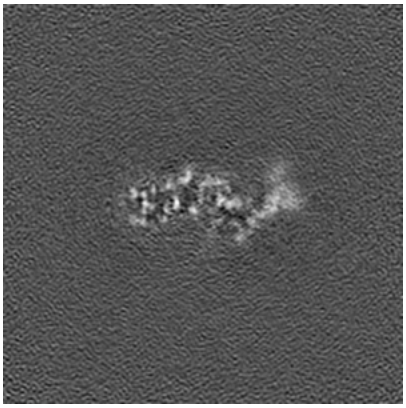

X Index: 100

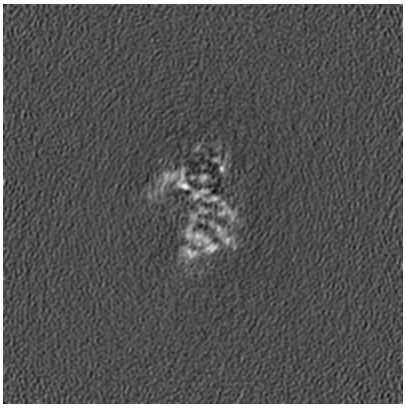

Y Index: 100


Z Index: 100

6.2.2 Raw map


X Index: 100

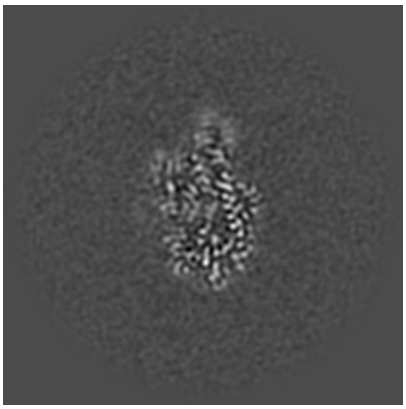

Y Index: 100

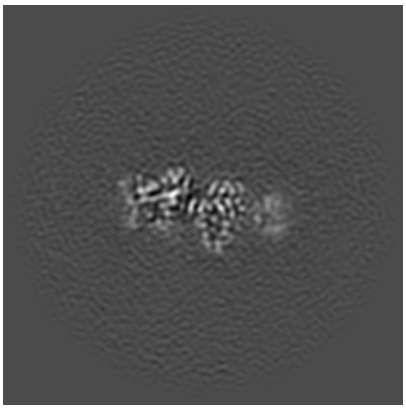

Z Index: 100

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

6.3 Largest variance slices (i)

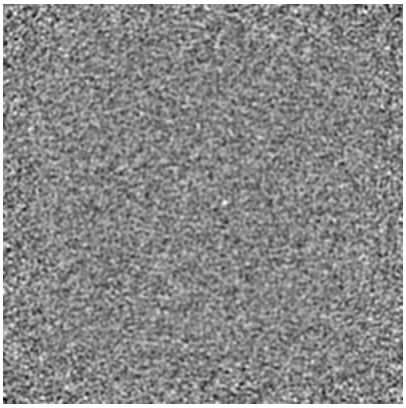
6.3.1 Primary map

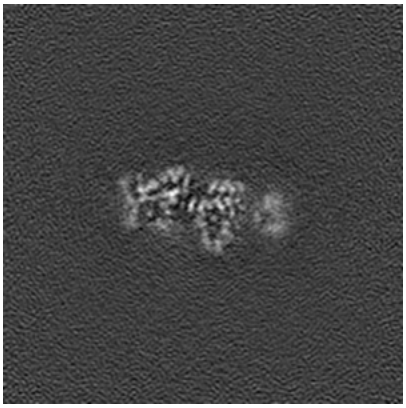

X Index: 99

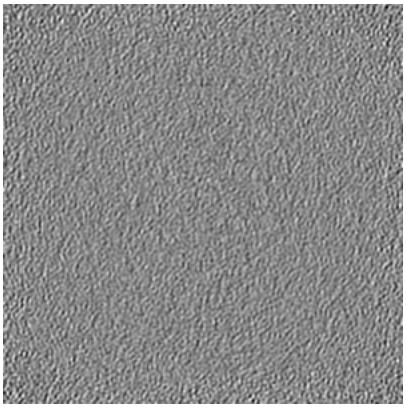

Y Index: 113

Z Index: 108

6.3.2 Raw map


X Index: 0

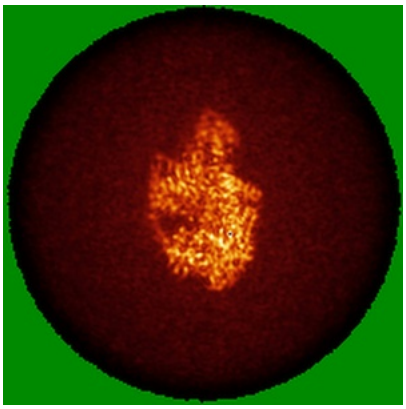

Y Index: 113

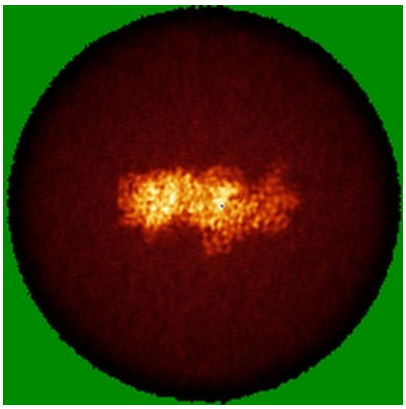

Z Index: 0

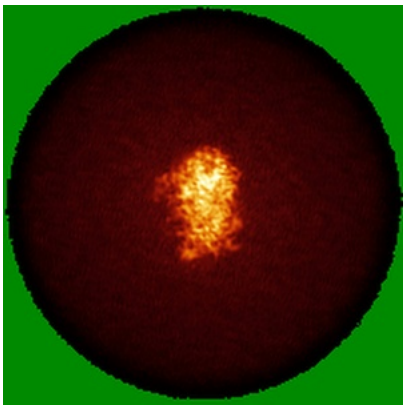
The images above show the largest variance slices of the map in three orthogonal directions.

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

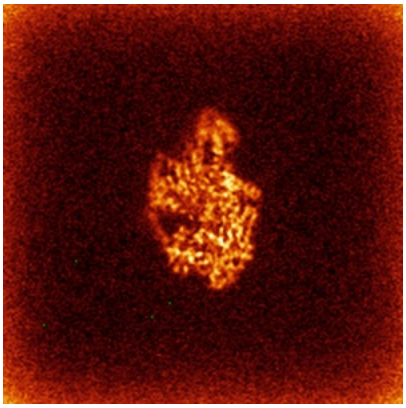
6.4.1 Primary map

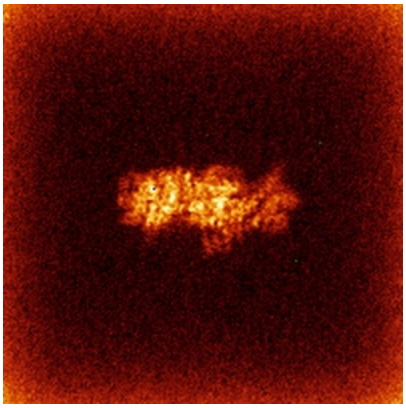

X

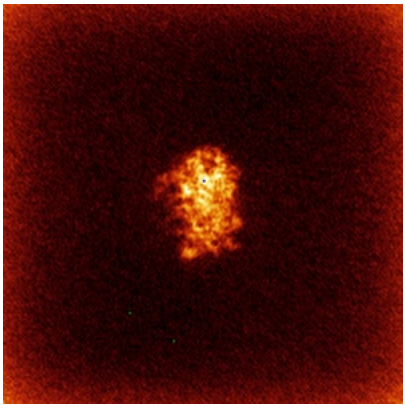

Y


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 0.08. 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.

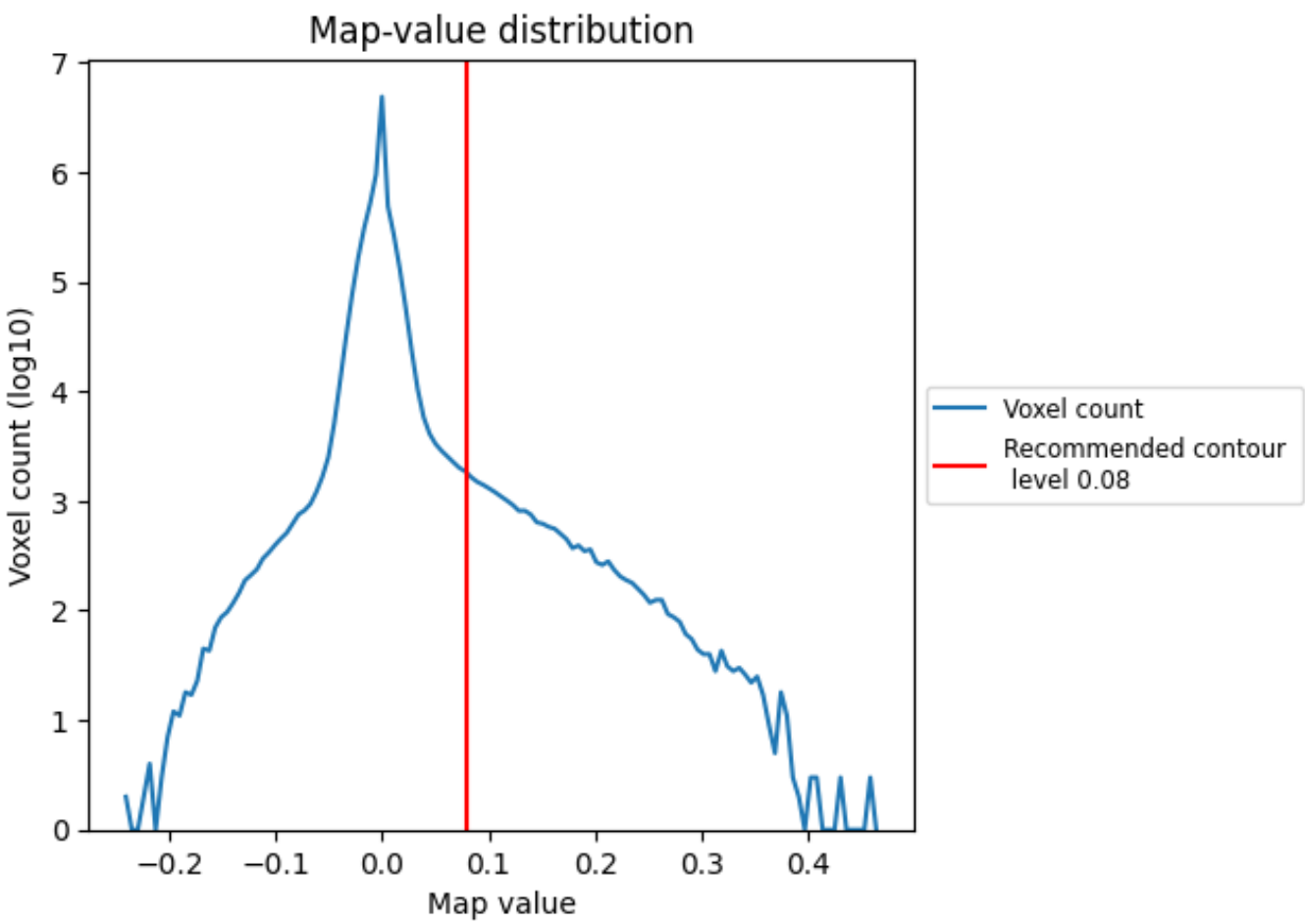
6.6 Mask visualisation (i)

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

7 Map analysis (i)

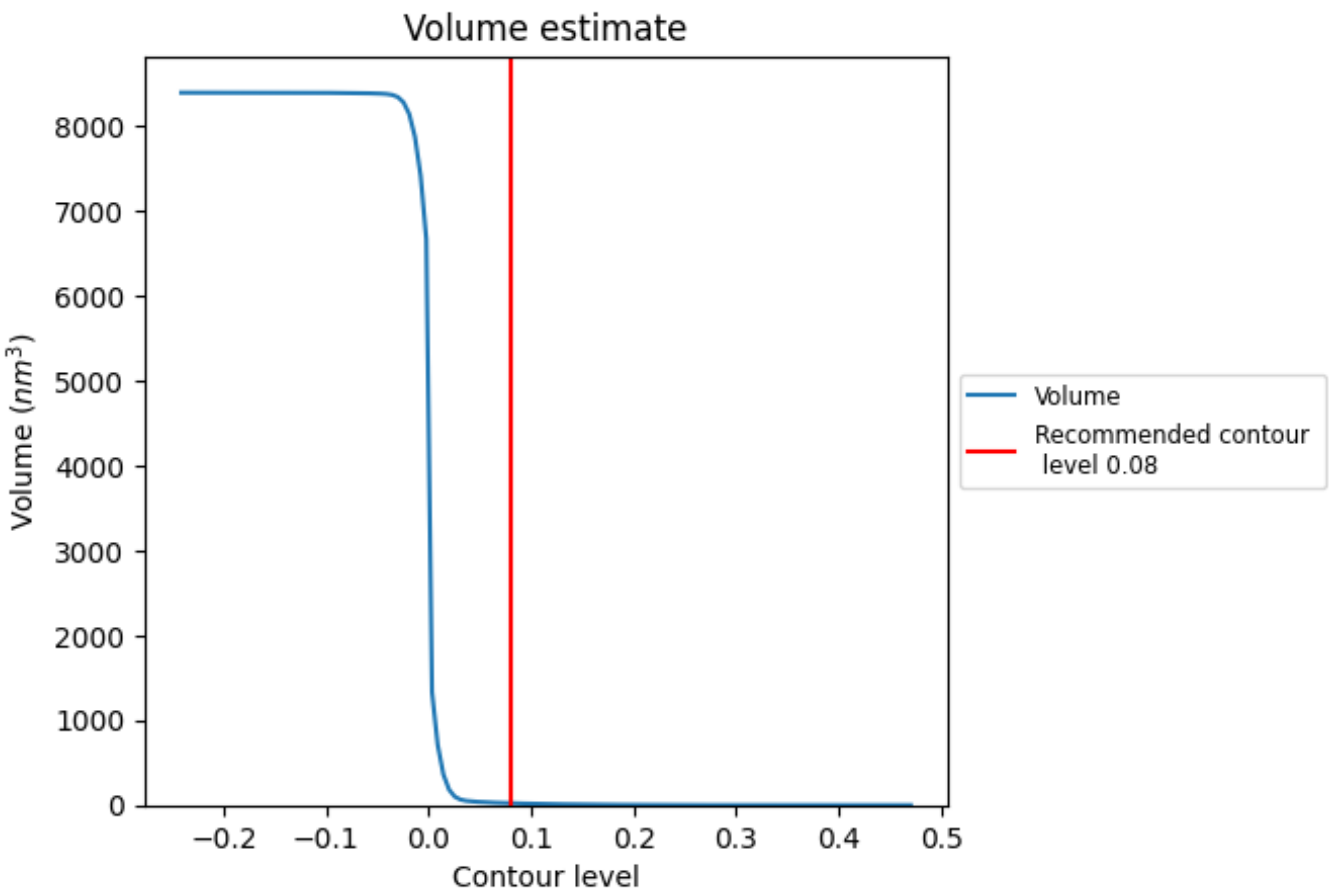
This section contains the results of statistical analysis of the map.

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.

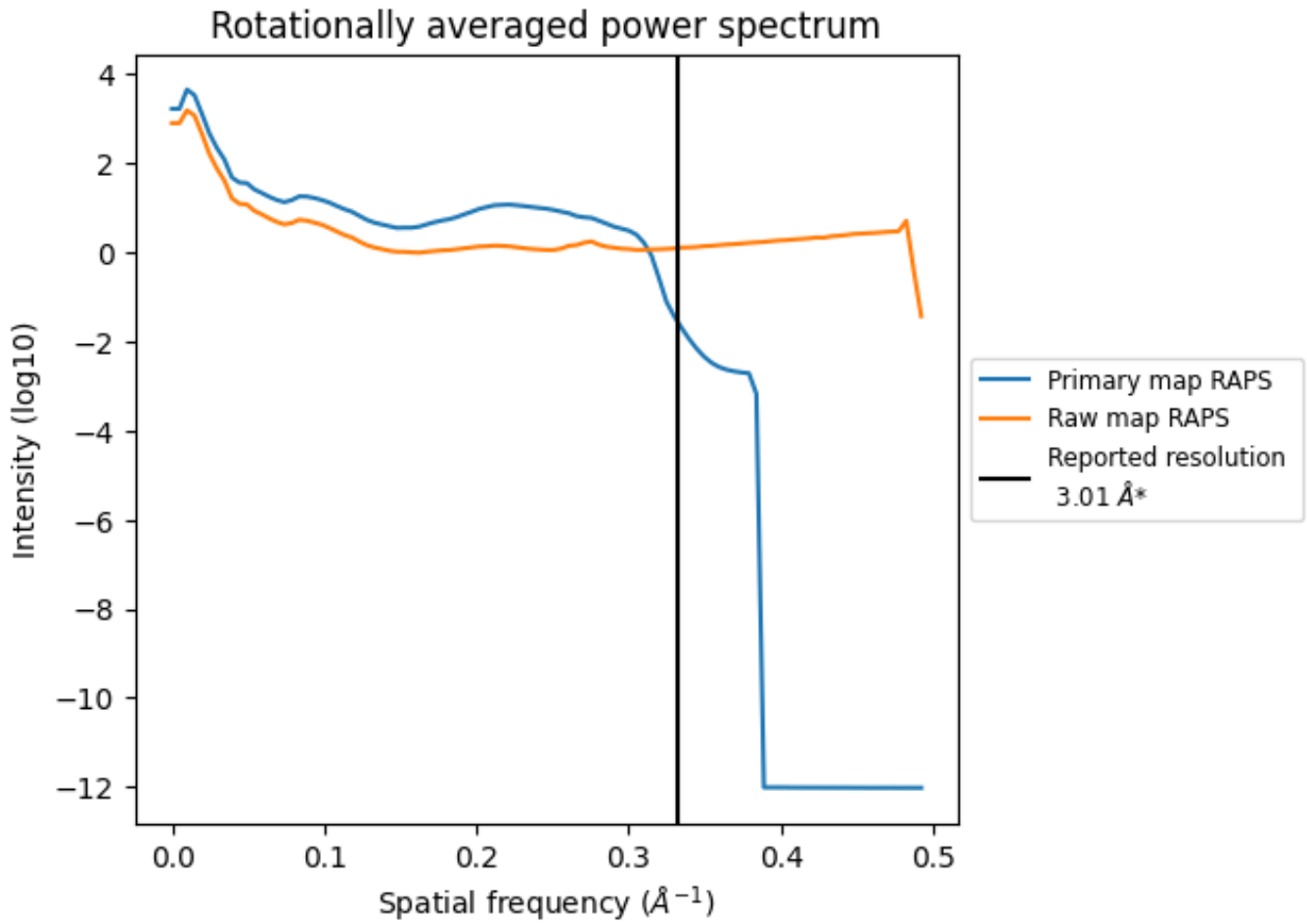
7.2 Volume estimate (i)



The volume at the recommended contour level is 23 nm^3; this corresponds to an approximate mass of 20 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)

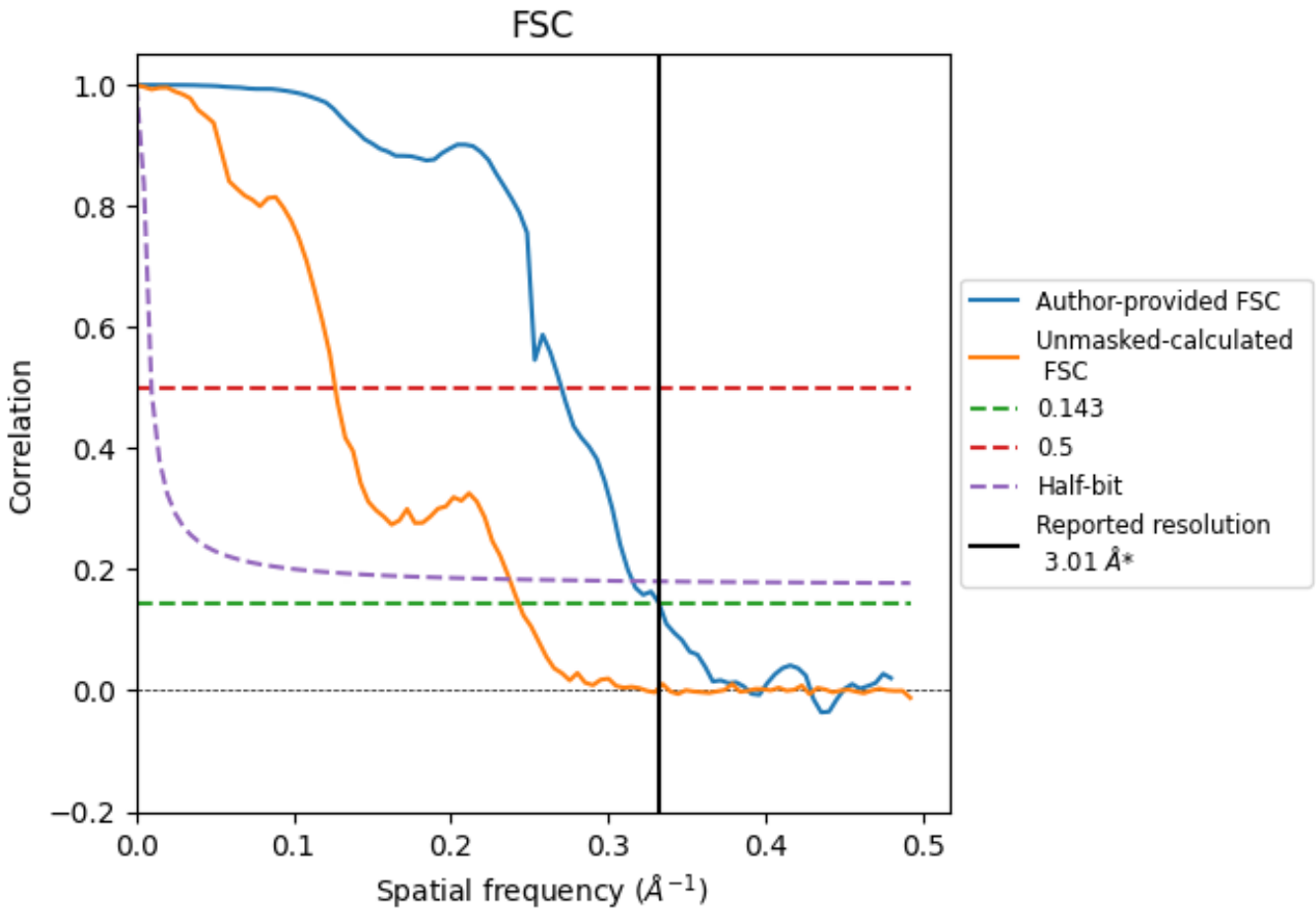


*Reported resolution corresponds to spatial frequency of 0.332 Å^{-1}

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 0.332 Å^{-1}

8.2 Resolution estimates (i)

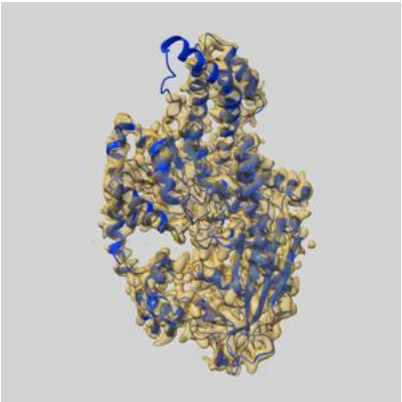
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.01	-	-
Author-provided FSC curve	3.01	3.70	3.17
Unmasked-calculated*	4.11	7.91	4.21

*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 4.11 differs from the reported value 3.01 by more than 10 %

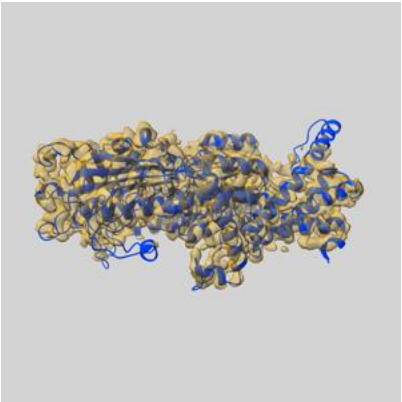
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-55193 and PDB model 9SSS. Per-residue inclusion information can be found in section 3 on page 4.

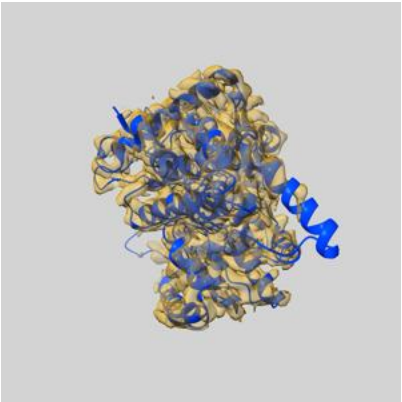
9.1 Map-model overlay (i)



X



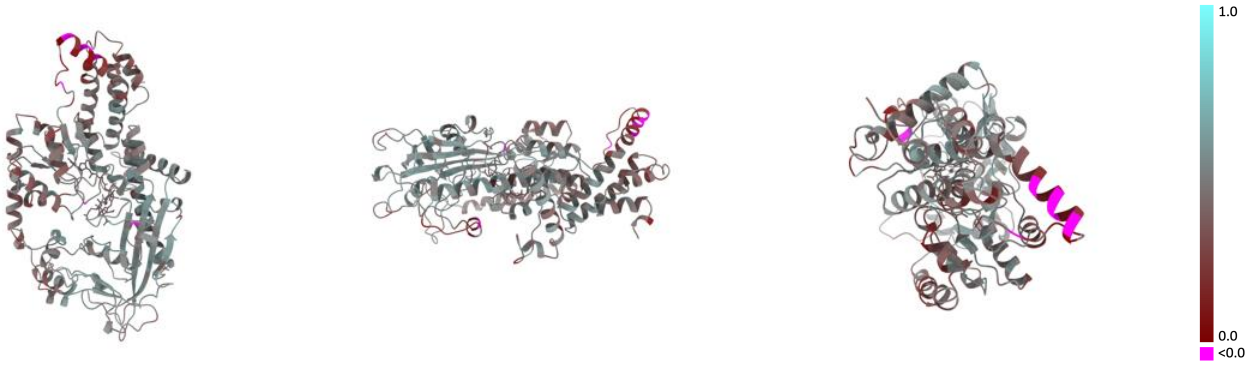
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.08 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.

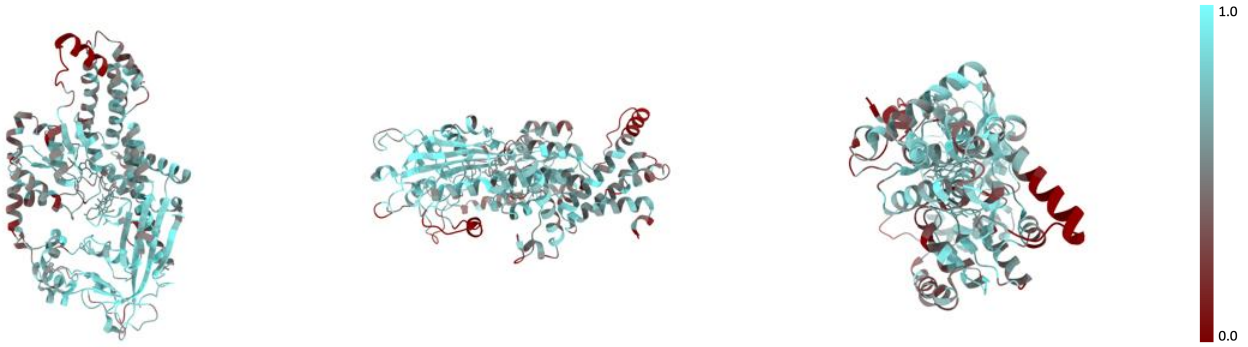
9.2 Q-score mapped to coordinate model (i)



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.

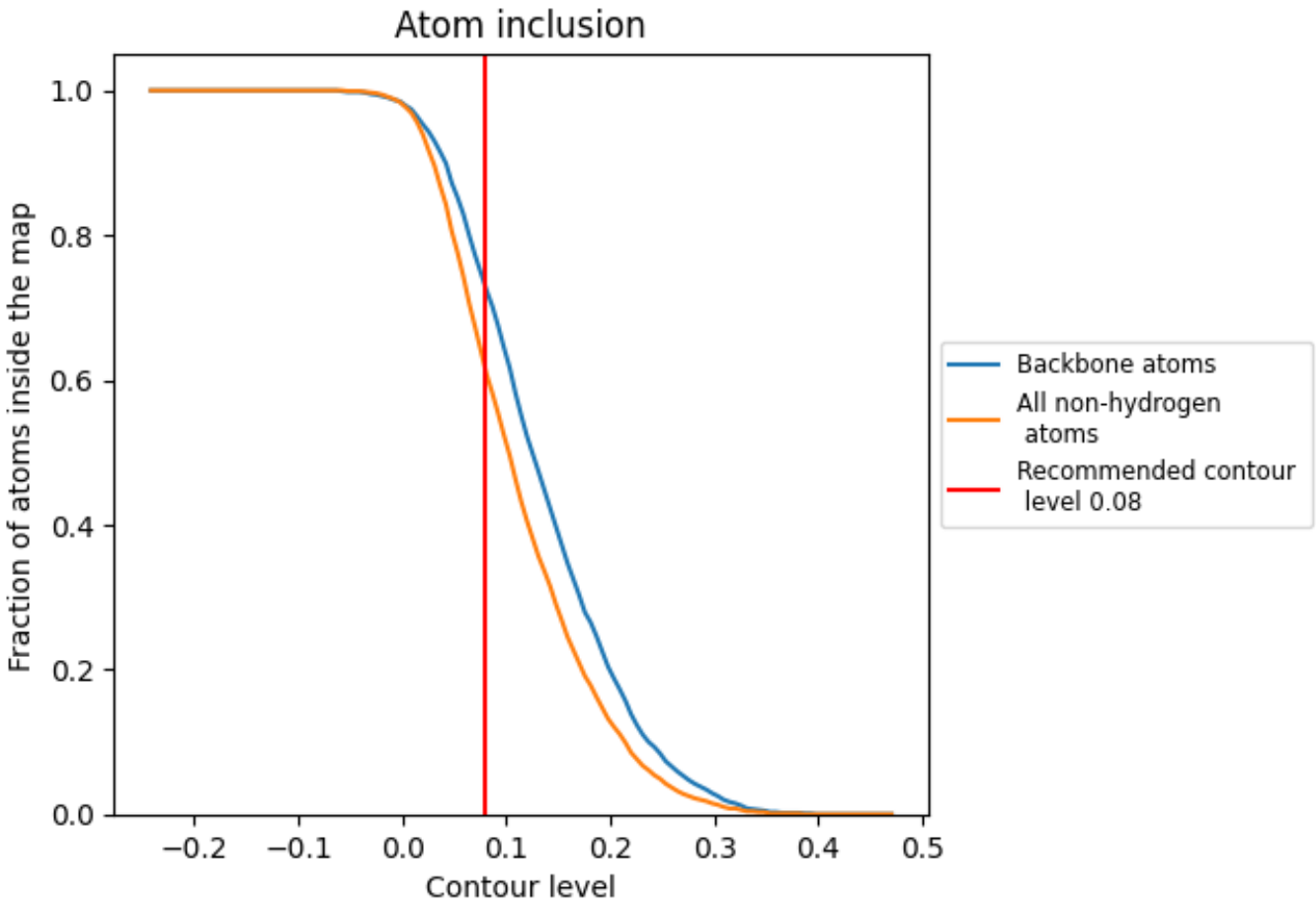
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.08).

9.4 Atom inclusion (i)



At the recommended contour level, 73% of all backbone atoms, 62% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.08) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6150	0.4410
A	0.6150	0.4410