



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:36 AM UTC

PDB ID : 6F6W / pdb_00006f6w
EMDB ID : EMD-4192
Title : Structure of Mycobacterium smegmatis RNA polymerase core
Authors : Kouba, T.; Barvik, I.; Krasny, L.
Deposited on : 2017-12-06
Resolution : 3.81 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:

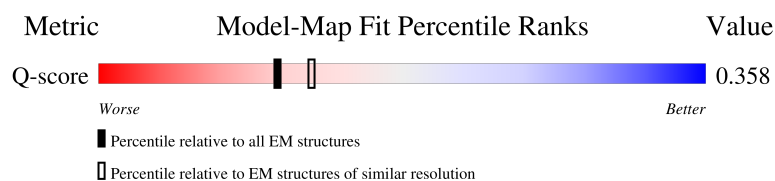
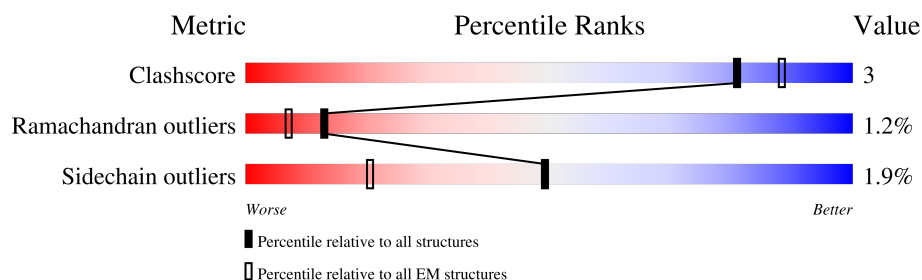
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.81 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	9193 (3.31 - 4.31)

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	350	
1	B	350	
2	C	1178	
3	D	1325	

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Mol	Chain	Length	Quality of chain	
			5%	29%
4	E	107	 A horizontal bar chart showing the quality of chain E. The bar is divided into three segments: a red segment at the beginning labeled '5%', a green segment in the middle labeled '66%', and a grey segment at the end labeled '29%'. The green segment is the largest, followed by the grey segment, and the red segment is the smallest.	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 21619 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	221	Total	C	N	O	S	0	0
			1648	1042	285	318	3		
1	B	233	Total	C	N	O	S	0	0
			1683	1062	295	324	2		

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	1090	Total	C	N	O	S	0	0
			8139	5093	1427	1585	34		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	VAL	-	expression tag	UNP P60281
C	1170	LEU	-	expression tag	UNP P60281
C	1171	ALA	-	expression tag	UNP P60281
C	1172	ARG	-	expression tag	UNP P60281
C	1173	HIS	-	expression tag	UNP P60281
C	1174	GLY	-	expression tag	UNP P60281
C	1175	GLY	-	expression tag	UNP P60281
C	1176	SER	-	expression tag	UNP P60281
C	1177	GLY	-	expression tag	UNP P60281
C	1178	ALA	-	expression tag	UNP P60281

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	1242	Total	C	N	O	S	0	0
			9554	5997	1721	1796	40		

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	VAL	-	expression tag	UNP A0QS66
D	1318	HIS	-	expression tag	UNP A0QS66
D	1319	HIS	-	expression tag	UNP A0QS66
D	1320	HIS	-	expression tag	UNP A0QS66
D	1321	HIS	-	expression tag	UNP A0QS66
D	1322	HIS	-	expression tag	UNP A0QS66
D	1323	HIS	-	expression tag	UNP A0QS66
D	1324	HIS	-	expression tag	UNP A0QS66
D	1325	HIS	-	expression tag	UNP A0QS66

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	E	76	Total	C	N	O	0	0
			592	378	100	114		

There is a discrepancy between the modelled and reference sequences:

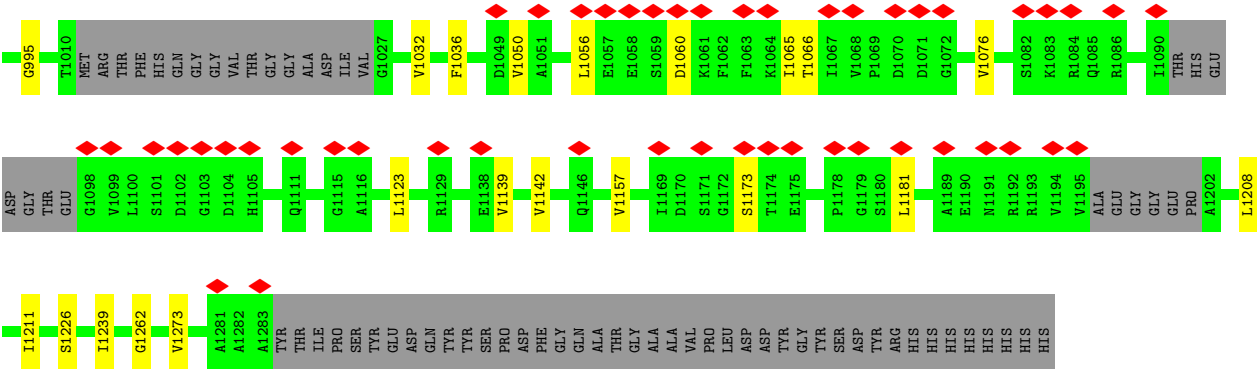
Chain	Residue	Modelled	Actual	Comment	Reference
E	1	VAL	-	expression tag	UNP A0QWT1

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

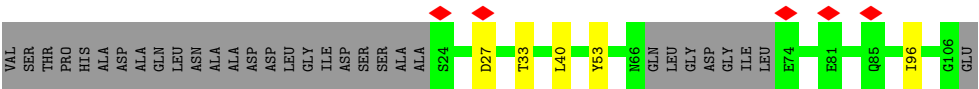
Mol	Chain	Residues	Atoms		AltConf
5	D	2	Total	Zn	0
			2	2	

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
6	D	1	Total	Mg	0
			1	1	



● Molecule 4: DNA-directed RNA polymerase subunit omega



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	163000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3800	Depositor
Magnification	78000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.753	Depositor
Minimum map value	-0.420	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.019	Depositor
Recommended contour level	0.1	Depositor
Map size ($\text{\AA}$)	278.72, 278.72, 278.72	wwPDB
Map dimensions	208, 208, 208	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.34, 1.34, 1.34	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: MG, ZN

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	1.09	0/1674	1.45	0/2281
1	B	1.12	0/1709	1.46	0/2335
2	C	1.07	0/8282	1.49	0/11265
3	D	1.08	0/9707	1.57	3/13140 (0.0%)
4	E	1.02	0/604	1.55	0/822
All	All	1.08	0/21976	1.52	3/29843 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	D	505	HIS	CA-C-N	5.14	127.42	120.38
3	D	505	HIS	C-N-CA	5.14	127.42	120.38
3	D	408	GLY	CA-C-O	-5.01	118.56	122.52

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	1648	0	1680	11	0
1	B	1683	0	1662	16	0
2	C	8139	0	7856	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	9554	0	9527	50	0
4	E	592	0	583	3	0
5	D	2	0	0	0	0
6	D	1	0	0	0	0
All	All	21619	0	21308	113	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 3.

All (113) 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)
2:C:839:ILE:HD11	2:C:865:ALA:HB2	1.65	0.78
3:D:883:VAL:HG11	3:D:1157:VAL:HG13	1.70	0.71
2:C:876:LEU:HD22	2:C:886:ILE:HD11	1.71	0.70
2:C:1126:VAL:HG12	3:D:12:ILE:HG22	1.74	0.70
2:C:649:ILE:HD13	2:C:693:ILE:HG22	1.77	0.67
2:C:47:VAL:HG11	2:C:491:LEU:HD22	1.78	0.64
2:C:717:TRP:HA	3:D:724:THR:HG22	1.78	0.64
3:D:456:VAL:HA	3:D:490:VAL:HG21	1.83	0.61
2:C:728:LEU:HB3	2:C:906:ILE:HG22	1.83	0.59
1:B:80:LEU:HD22	1:B:138:LEU:HD21	1.83	0.59
2:C:45:LEU:HD11	2:C:443:LYS:HD3	1.86	0.56
1:A:80:LEU:HD22	1:A:138:LEU:HD21	1.88	0.56
3:D:699:LEU:HD13	3:D:711:THR:HG21	1.88	0.55
2:C:924:GLU:OE2	2:C:1020:TYR:N	2.39	0.55
3:D:488:GLU:HA	3:D:491:ILE:HG22	1.88	0.54
2:C:641:ILE:HD11	2:C:687:VAL:HG13	1.89	0.53
2:C:1059:PHE:CZ	2:C:1079:LEU:HD13	2.44	0.53
2:C:578:VAL:HG22	2:C:582:THR:HB	1.90	0.53
3:D:432:VAL:HG22	3:D:536:PHE:CE2	2.44	0.53
1:A:34:LEU:HD11	1:B:218:LEU:HD21	1.89	0.52
3:D:1050:VAL:HG12	3:D:1076:VAL:HG23	1.92	0.52
3:D:589:THR:HG21	3:D:687:MET:CG	2.40	0.52
2:C:166:VAL:HB	2:C:372:VAL:HG22	1.93	0.51
1:A:83:LEU:HD21	1:A:107:ALA:HB1	1.92	0.51
3:D:995:GLY:HA2	3:D:1157:VAL:HG11	1.92	0.51
3:D:1032:VAL:HG22	3:D:1142:VAL:HG21	1.92	0.50
3:D:579:LEU:O	3:D:806:ALA:HB1	2.11	0.50
2:C:932:LYS:HG2	2:C:1018:TYR:CE2	2.46	0.50
3:D:723:ALA:O	3:D:726:SER:OG	2.29	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:849:PHE:O	3:D:852:THR:OG1	2.29	0.50
2:C:911:HIS:ND1	3:D:801:ILE:HG12	2.27	0.50
2:C:602:MET:HG2	2:C:1024:LEU:HD11	1.93	0.49
3:D:298:VAL:HG23	3:D:1239:ILE:HD12	1.95	0.49
3:D:661:TRP:CZ3	3:D:663:ALA:HB2	2.48	0.49
1:A:51:VAL:HG11	1:A:138:LEU:HD23	1.95	0.48
2:C:176:VAL:HG12	2:C:195:VAL:HG22	1.95	0.48
3:D:839:PHE:CE1	3:D:843:LEU:HD11	2.47	0.48
2:C:607:VAL:HG22	2:C:608:PRO:HD2	1.95	0.48
3:D:589:THR:HG21	3:D:687:MET:HG3	1.95	0.48
2:C:302:VAL:HG12	2:C:500:PHE:HB3	1.96	0.47
1:A:82:GLY:HA3	1:A:123:MET:HE1	1.97	0.47
2:C:215:VAL:HG23	2:C:225:VAL:HA	1.96	0.47
3:D:622:MET:SD	3:D:629:VAL:HG22	2.53	0.47
2:C:913:VAL:HB	2:C:914:PRO:HD3	1.95	0.47
2:C:588:LEU:HD11	2:C:596:ALA:HB1	1.96	0.47
2:C:45:LEU:HD12	2:C:155:GLY:HA3	1.97	0.47
2:C:1065:TRP:CH2	3:D:874:ARG:HA	2.50	0.46
4:E:40:LEU:HD22	4:E:96:ILE:HG23	1.97	0.46
1:A:40:ARG:HB3	1:B:33:THR:HG22	1.98	0.46
1:B:21:PHE:HB2	1:B:194:ILE:HG23	1.98	0.46
2:C:943:VAL:HG22	2:C:944:PRO:HD2	1.97	0.46
3:D:532:PHE:CZ	3:D:541:MET:HE1	2.50	0.46
3:D:1123:LEU:HD22	3:D:1208:LEU:HB2	1.97	0.45
3:D:990:ILE:HG22	3:D:990:ILE:O	2.16	0.45
1:B:22:VAL:HG12	1:B:193:ILE:HD13	1.99	0.45
2:C:224:PRO:O	2:C:227:VAL:HG22	2.17	0.45
1:B:23:ILE:HD12	1:B:192:LEU:HD23	1.99	0.45
2:C:369:LEU:HD22	2:C:503:ILE:HD11	1.99	0.45
2:C:466:VAL:HG23	3:D:853:HIS:CE1	2.52	0.45
1:A:34:LEU:HD11	1:B:218:LEU:HD11	1.98	0.44
1:B:34:LEU:HD11	1:B:192:LEU:HD22	1.99	0.44
2:C:876:LEU:HD23	2:C:1021:ILE:HD13	1.99	0.44
3:D:97:LEU:HD11	3:D:317:VAL:HG23	1.98	0.44
3:D:413:PHE:HB3	3:D:1226:SER:HA	1.99	0.44
3:D:951:LEU:HD22	3:D:956:ILE:HD11	1.98	0.44
3:D:525:HIS:CG	3:D:526:PRO:HD2	2.53	0.44
2:C:607:VAL:HG22	2:C:608:PRO:CD	2.48	0.44
3:D:1273:VAL:HG21	4:E:53:TYR:HA	2.00	0.44
3:D:579:LEU:HD23	3:D:807:THR:HB	2.00	0.43
3:D:823:VAL:HG11	3:D:851:ASN:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:494:HIS:CE1	3:D:545:LEU:HD11	2.52	0.43
1:A:218:LEU:HD21	1:B:34:LEU:CD2	2.48	0.43
2:C:583:ALA:HB1	2:C:621:MET:HG3	2.01	0.43
1:B:172:LEU:HD21	1:B:199:LYS:HD3	2.01	0.43
2:C:77:LEU:HA	2:C:80:VAL:HG12	2.00	0.43
1:A:146:TYR:CG	2:C:734:GLU:HG2	2.54	0.43
2:C:412:ARG:N	2:C:413:PRO:HD2	2.34	0.43
4:E:53:TYR:CZ	4:E:96:ILE:HD13	2.54	0.43
1:B:171:VAL:HA	1:B:198:THR:HA	1.99	0.43
2:C:454:LEU:HD13	2:C:459:ALA:HB2	2.01	0.43
1:B:22:VAL:HG12	1:B:193:ILE:CD1	2.49	0.42
2:C:94:MET:HE2	2:C:400:VAL:HG23	2.01	0.42
3:D:1056:LEU:HD13	3:D:1065:ILE:HG13	2.01	0.42
3:D:491:ILE:HD11	3:D:514:PRO:HB2	2.02	0.42
2:C:364:PHE:CE1	2:C:503:ILE:HD12	2.54	0.42
1:B:78:LEU:HD11	3:D:611:VAL:HB	2.01	0.42
1:B:107:ALA:O	1:B:110:ILE:HG22	2.19	0.42
3:D:812:GLN:HA	3:D:815:THR:HG22	2.00	0.42
2:C:493:VAL:HG13	2:C:494:TYR:N	2.34	0.42
2:C:281:GLU:HA	2:C:284:GLN:HE21	1.84	0.42
2:C:885:VAL:HG11	3:D:429:VAL:CG1	2.49	0.42
3:D:901:ALA:HB1	3:D:910:ILE:HG23	2.01	0.42
3:D:589:THR:HG21	3:D:687:MET:HG2	2.02	0.41
3:D:746:GLU:O	3:D:749:GLU:HB2	2.20	0.41
3:D:934:ASN:O	3:D:935:VAL:C	2.63	0.41
2:C:711:LEU:HD12	2:C:1017:GLY:O	2.21	0.41
1:A:44:SER:OG	1:A:45:SER:N	2.53	0.41
1:B:38:LEU:HD23	1:B:38:LEU:HA	1.97	0.41
3:D:277:LEU:HA	3:D:280:VAL:HG12	2.03	0.41
2:C:454:LEU:HD12	2:C:454:LEU:O	2.20	0.41
2:C:891:PRO:O	2:C:893:GLU:N	2.54	0.41
2:C:711:LEU:HD13	2:C:1018:TYR:CE2	2.56	0.41
2:C:728:LEU:HB2	2:C:732:LEU:HD22	2.03	0.41
3:D:14:LEU:HD22	3:D:106:TYR:CZ	2.56	0.41
1:B:34:LEU:HD12	1:B:38:LEU:HD12	2.02	0.41
2:C:1065:TRP:CD1	3:D:877:VAL:HG11	2.56	0.40
3:D:444:PRO:HD2	3:D:447:MET:HE2	2.03	0.40
3:D:876:LEU:HD12	3:D:1211:ILE:CD1	2.51	0.40
3:D:453:LYS:HB3	3:D:454:PRO:HD3	2.03	0.40
3:D:655:TRP:CE3	3:D:655:TRP:HA	2.55	0.40
3:D:1036:PHE:CE2	3:D:1139:VAL:HG21	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:ASN:HB2	1:A:157:ALA:HB3	2.03	0.40
2:C:710:LEU:HD13	2:C:732:LEU:HD21	2.03	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	219/350 (63%)	199 (91%)	15 (7%)	5 (2%)	5	30
1	B	231/350 (66%)	195 (84%)	32 (14%)	4 (2%)	7	34
2	C	1084/1178 (92%)	966 (89%)	105 (10%)	13 (1%)	10	40
3	D	1230/1325 (93%)	1114 (91%)	105 (8%)	11 (1%)	14	46
4	E	72/107 (67%)	65 (90%)	6 (8%)	1 (1%)	9	37
All	All	2836/3310 (86%)	2539 (90%)	263 (9%)	34 (1%)	13	40

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	181	THR
2	C	940	ALA
3	D	935	VAL
1	A	149	ALA
1	A	187	THR
1	A	200	ASN
2	C	572	VAL
2	C	892	VAL
2	C	938	ASP
3	D	238	GLU
3	D	1060	ASP
3	D	1262	GLY

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Mol	Chain	Res	Type
2	C	22	SER
2	C	789	ASP
2	C	937	ILE
3	D	492	ALA
3	D	625	GLY
1	B	7	PRO
1	B	64	THR
2	C	525	ASP
2	C	913	VAL
2	C	981	GLY
3	D	91	ARG
3	D	653	ASN
3	D	1173	SER
4	E	27	ASP
1	B	28	PRO
1	B	83	LEU
3	D	900	LEU
1	A	4	SER
2	C	1107	GLY
3	D	990	ILE
2	C	543	GLY
2	C	887	GLY

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	181/297 (61%)	174 (96%)	7 (4%)	28	52
1	B	174/297 (59%)	170 (98%)	4 (2%)	44	63
2	C	844/988 (85%)	823 (98%)	21 (2%)	42	61
3	D	992/1103 (90%)	983 (99%)	9 (1%)	70	74
4	E	62/86 (72%)	61 (98%)	1 (2%)	55	68
All	All	2253/2771 (81%)	2211 (98%)	42 (2%)	49	66

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	SER
1	A	84	VAL
1	A	127	THR
1	A	166	SER
1	A	188	ASP
1	A	194	ILE
1	A	195	ASP
1	B	105	VAL
1	B	106	THR
1	B	111	VAL
1	B	196	VAL
2	C	156	THR
2	C	207	VAL
2	C	313	LEU
2	C	505	THR
2	C	572	VAL
2	C	573	SER
2	C	578	VAL
2	C	592	ASP
2	C	632	VAL
2	C	706	LEU
2	C	727	ILE
2	C	728	LEU
2	C	733	VAL
2	C	749	ASP
2	C	870	ILE
2	C	907	ILE
2	C	943	VAL
2	C	965	THR
2	C	976	LEU
2	C	1041	SER
2	C	1078	GLU
3	D	83	THR
3	D	277	LEU
3	D	432	VAL
3	D	580	ASP
3	D	655	TRP
3	D	656	LYS
3	D	992	GLU
3	D	1066	THR
3	D	1181	LEU
4	E	33	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	151	GLN
2	C	160	ASN
2	C	236	ASN
2	C	284	GLN
2	C	366	ASN
2	C	433	GLN
2	C	434	ASN
2	C	467	HIS
2	C	484	ASN
2	C	576	GLN
2	C	676	ASN
3	D	267	ASN
3	D	303	GLN
3	D	341	ASN
3	D	368	ASN
3	D	505	HIS
3	D	609	GLN
3	D	697	ASN
3	D	710	GLN
3	D	809	ASN
3	D	851	ASN
3	D	1033	GLN
3	D	1085	GLN
3	D	1161	GLN
3	D	1240	ASN
4	E	66	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

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.

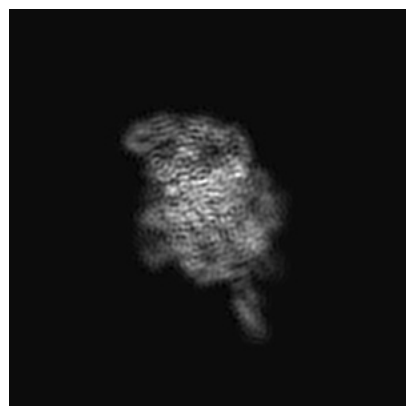
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4192. These allow visual inspection of the internal detail of the map and identification of artifacts.

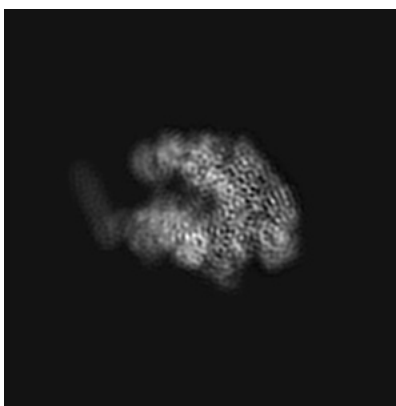
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

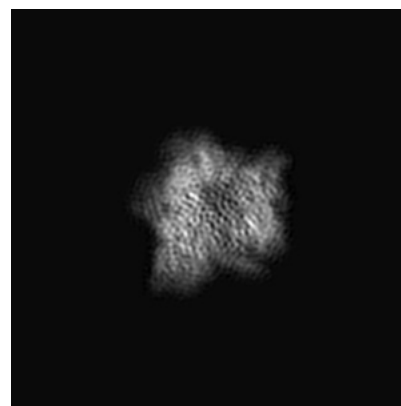
6.1.1 Primary map



X

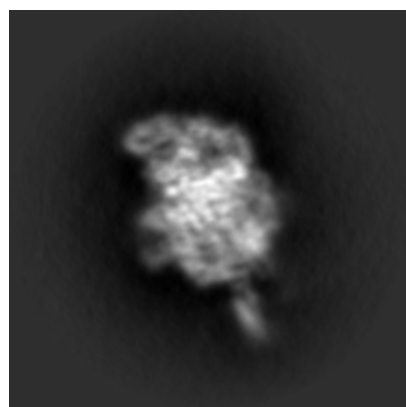


Y

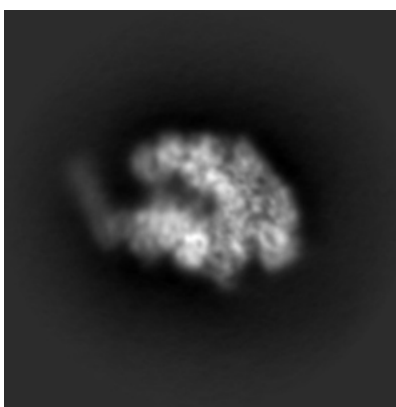


Z

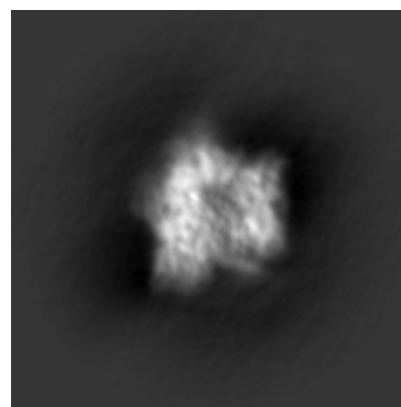
6.1.2 Raw map



X



Y

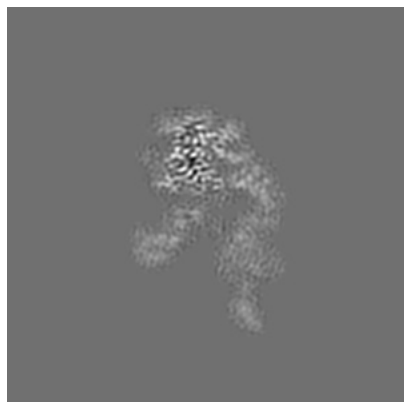


Z

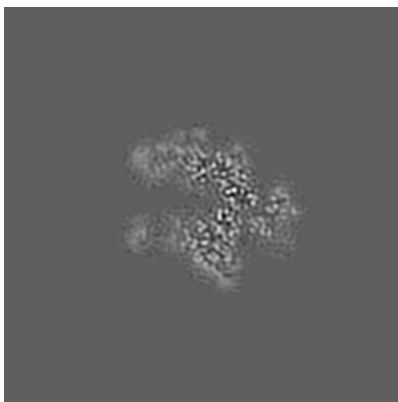
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

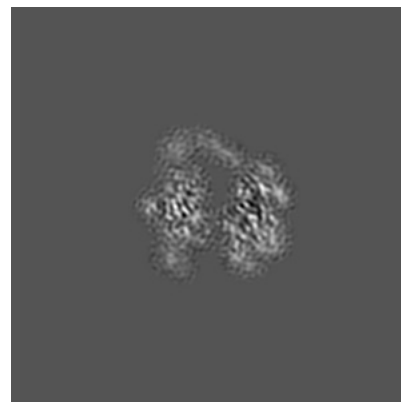
6.2.1 Primary map



X Index: 104

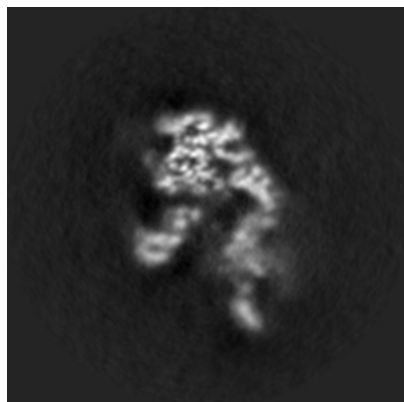


Y Index: 104

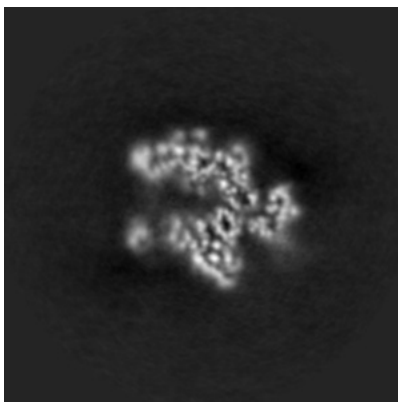


Z Index: 104

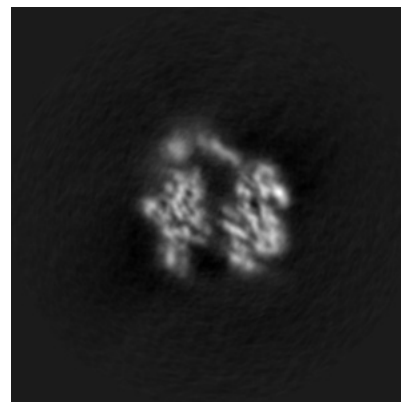
6.2.2 Raw map



X Index: 104



Y Index: 104

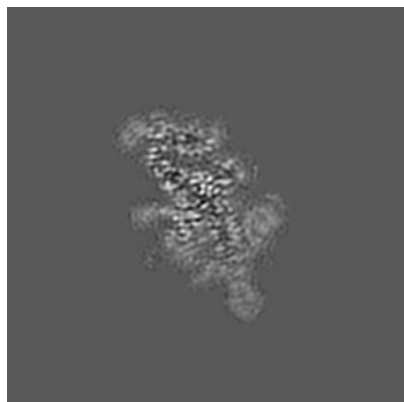


Z Index: 104

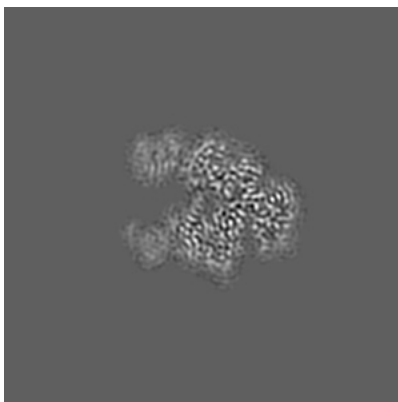
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

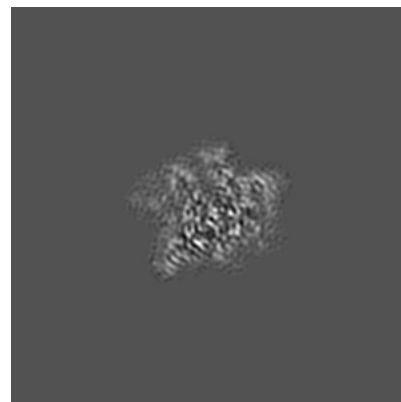
6.3.1 Primary map



X Index: 90



Y Index: 97

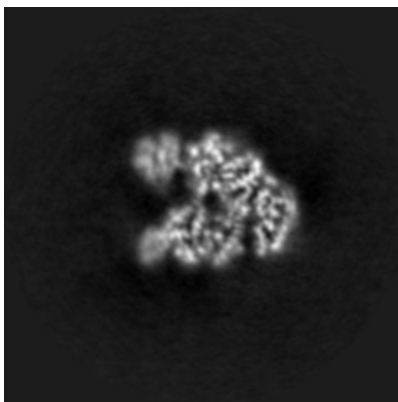


Z Index: 121

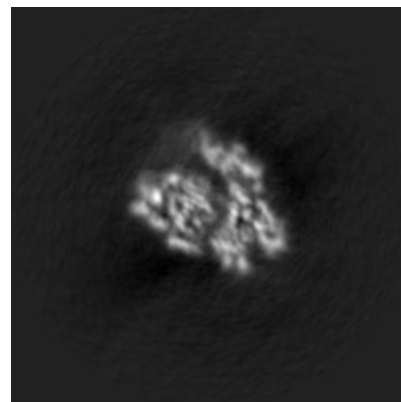
6.3.2 Raw map



X Index: 90



Y Index: 93

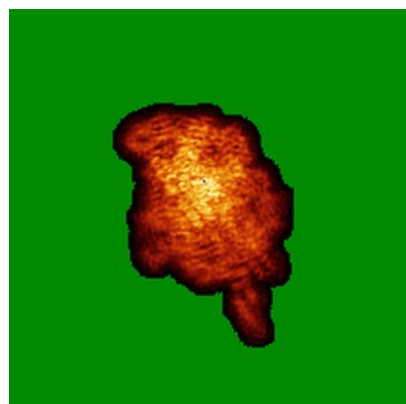


Z Index: 112

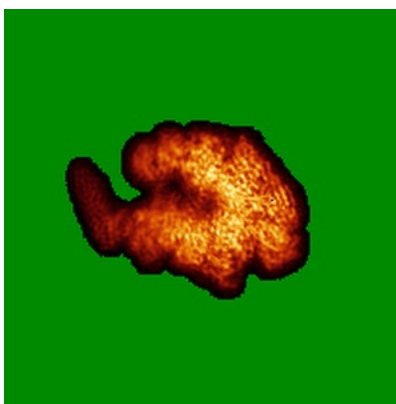
The images above show the largest variance slices of the map in three orthogonal directions.

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

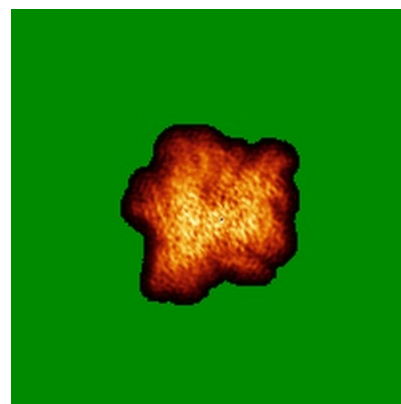
6.4.1 Primary map



X

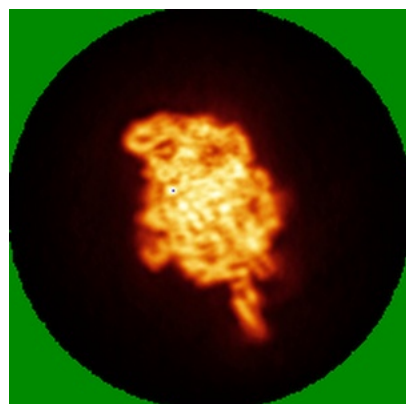


Y

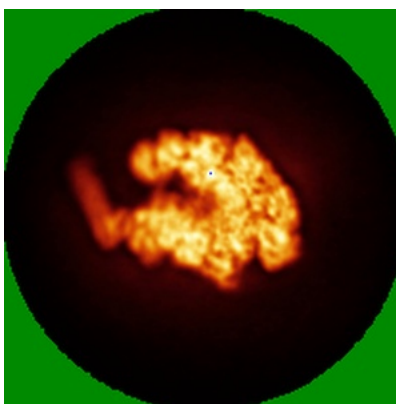


Z

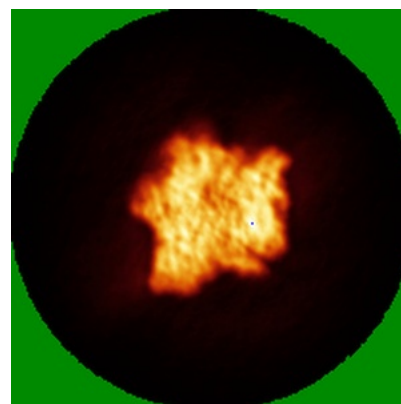
6.4.2 Raw map



X



Y

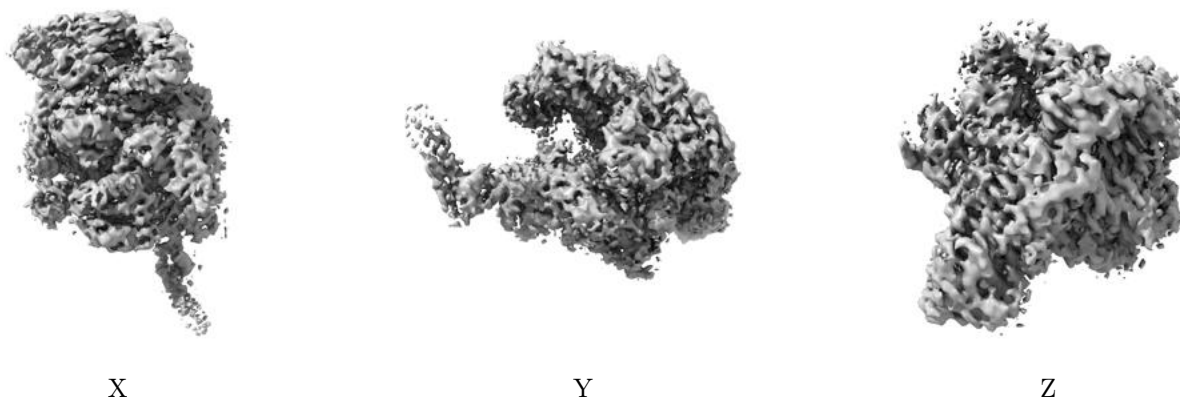


Z

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

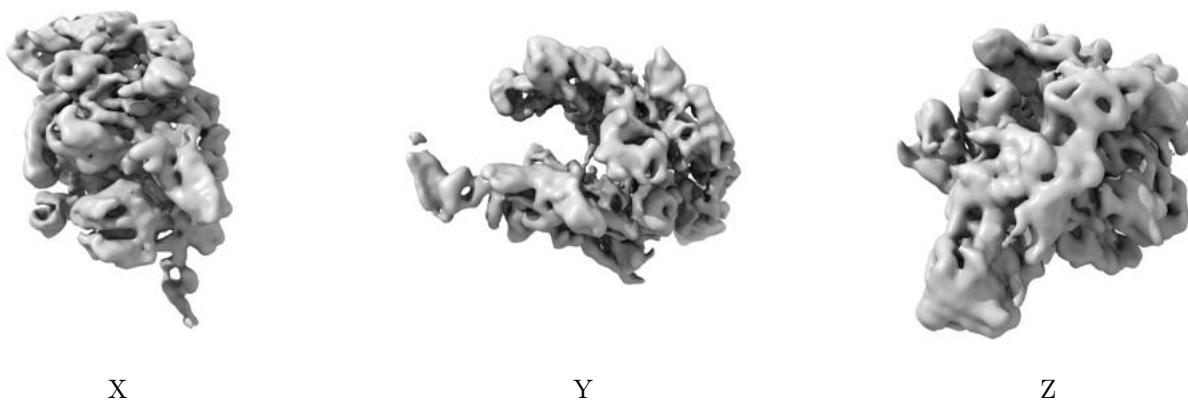
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

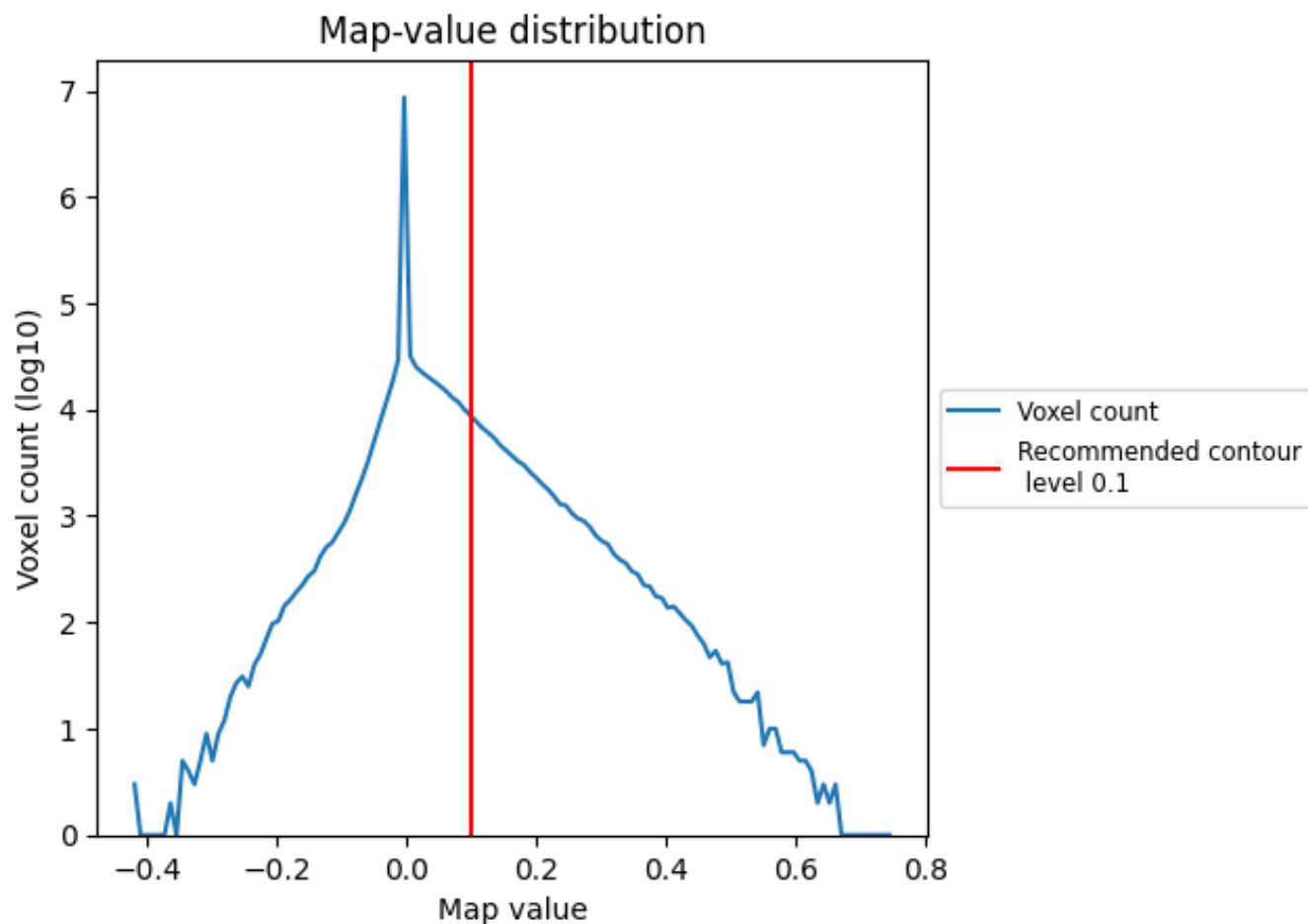
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

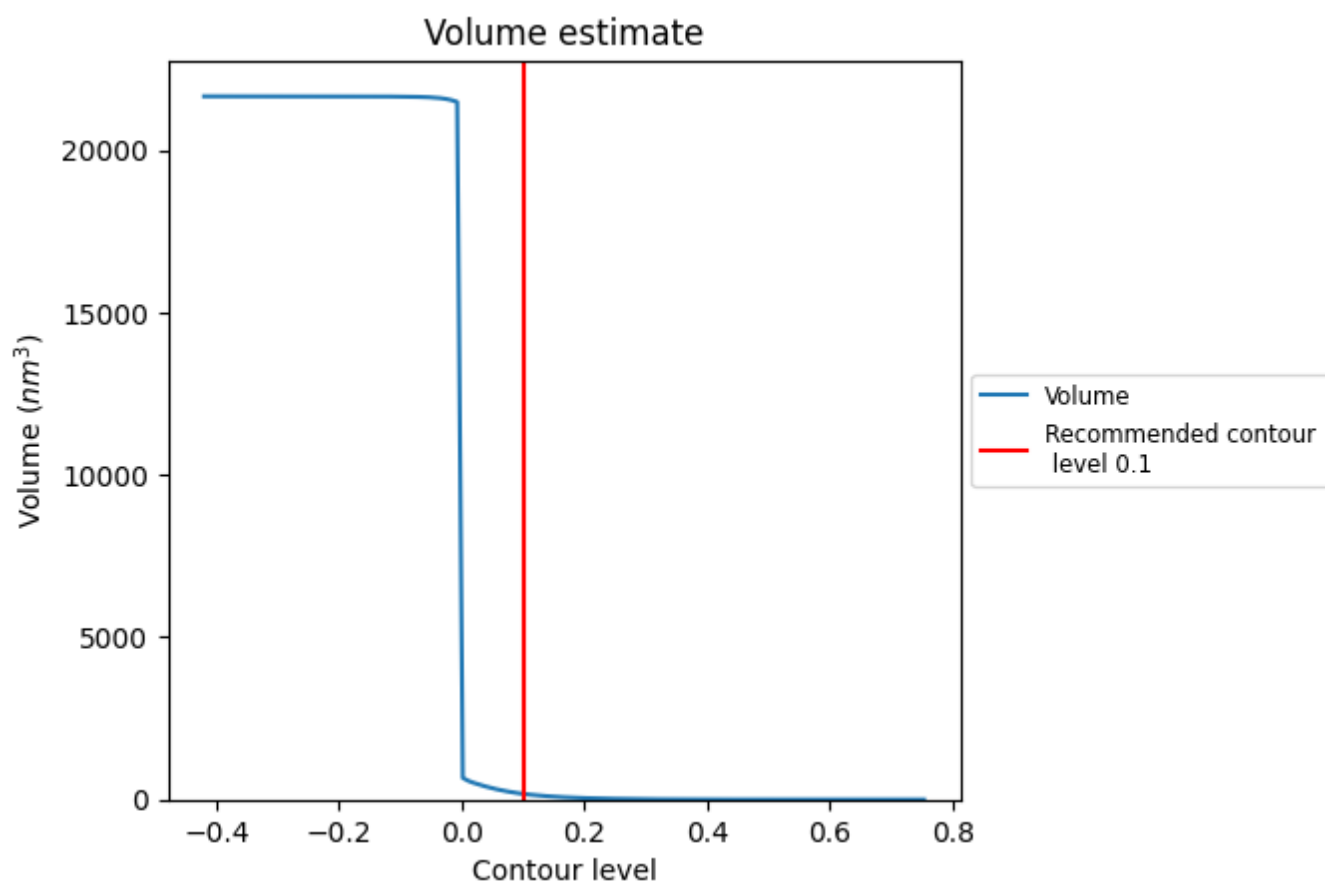
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

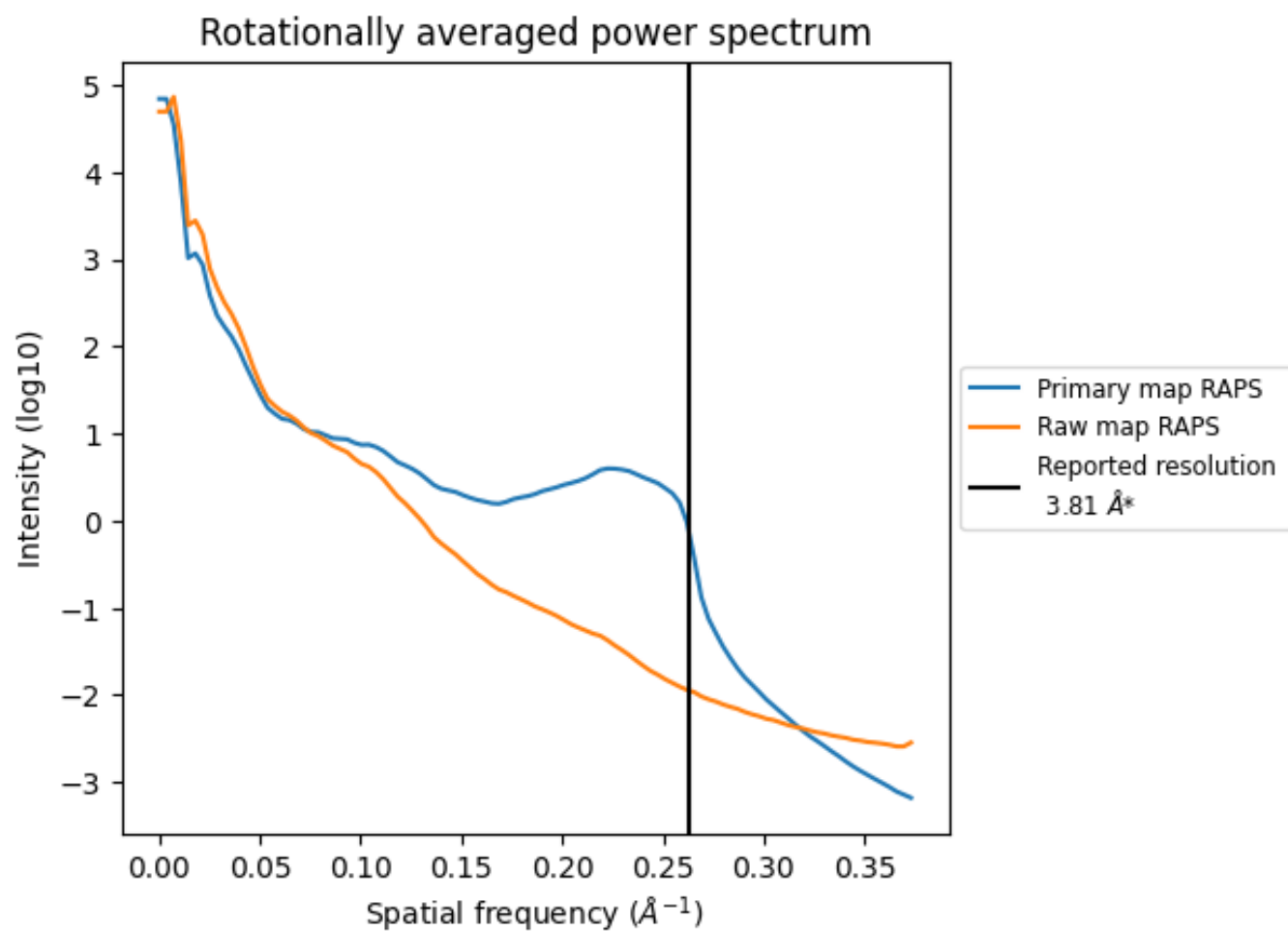
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 176 nm³; this corresponds to an approximate mass of 159 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 ⓘ

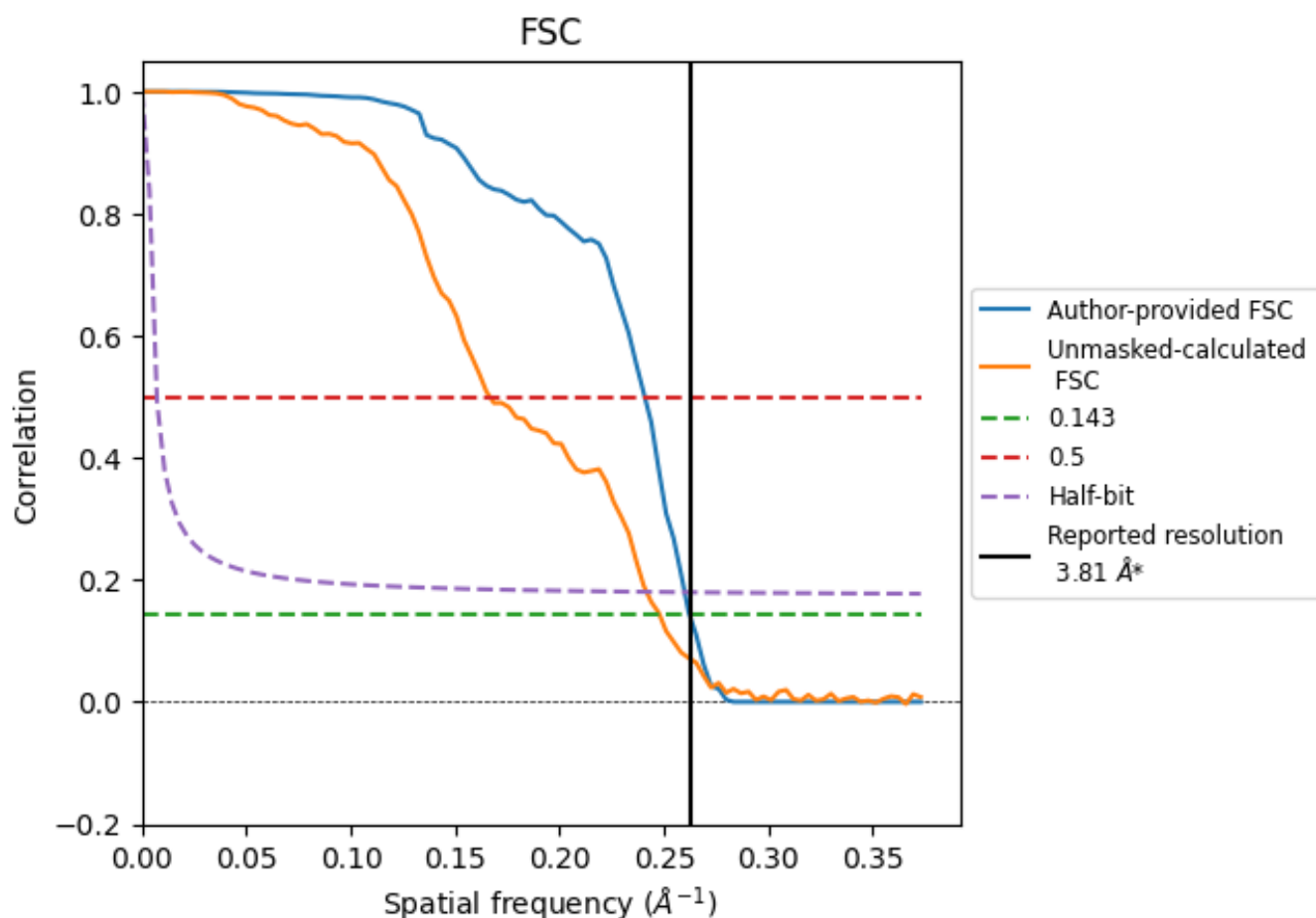


*Reported resolution corresponds to spatial frequency of 0.262 Å⁻¹

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 0.262 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

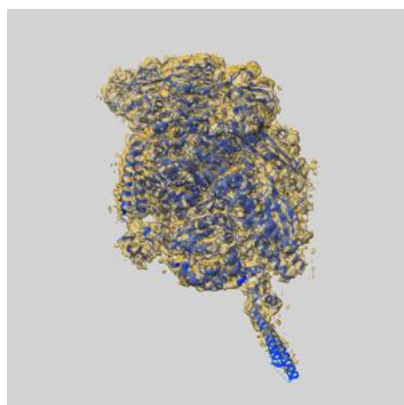
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.81	-	-
Author-provided FSC curve	3.81	4.15	3.85
Unmasked-calculated*	4.03	6.00	4.14

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

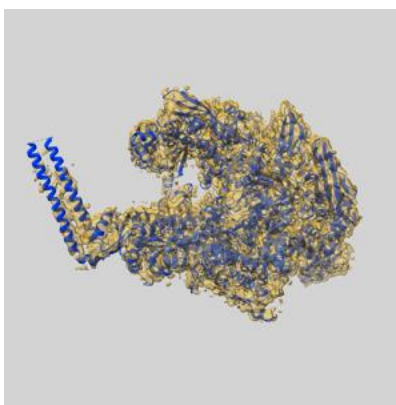
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-4192 and PDB model 6F6W. Per-residue inclusion information can be found in section [3](#) on page [6](#).

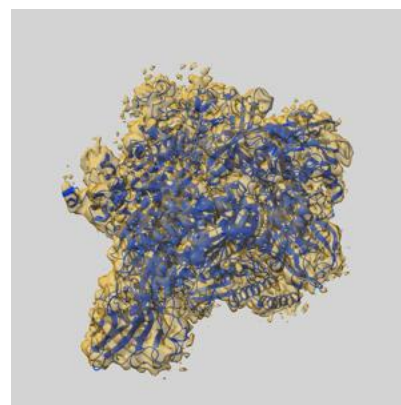
9.1 Map-model overlay [i](#)



X



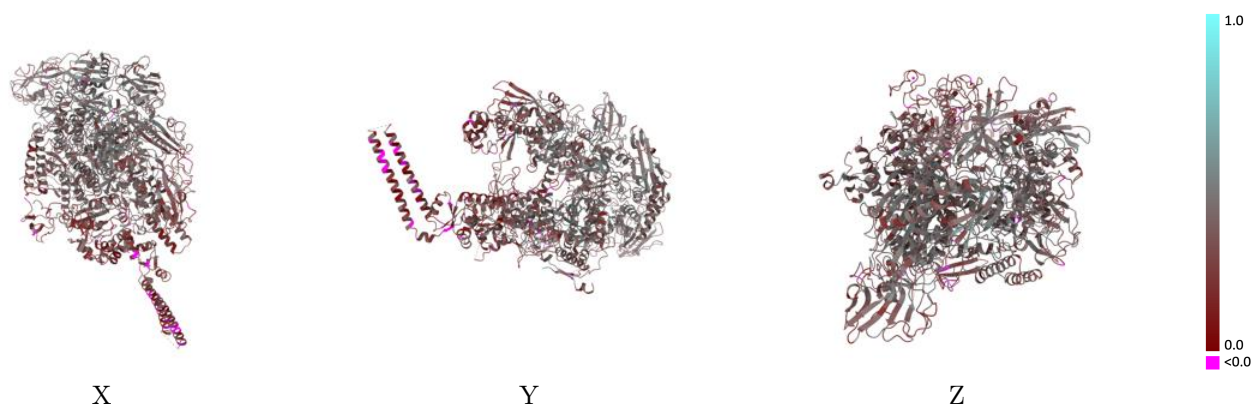
Y



Z

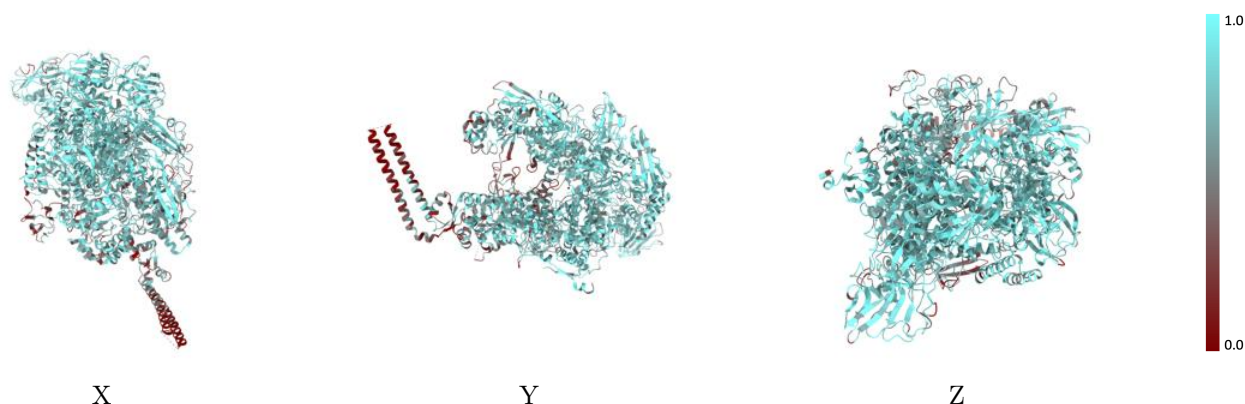
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



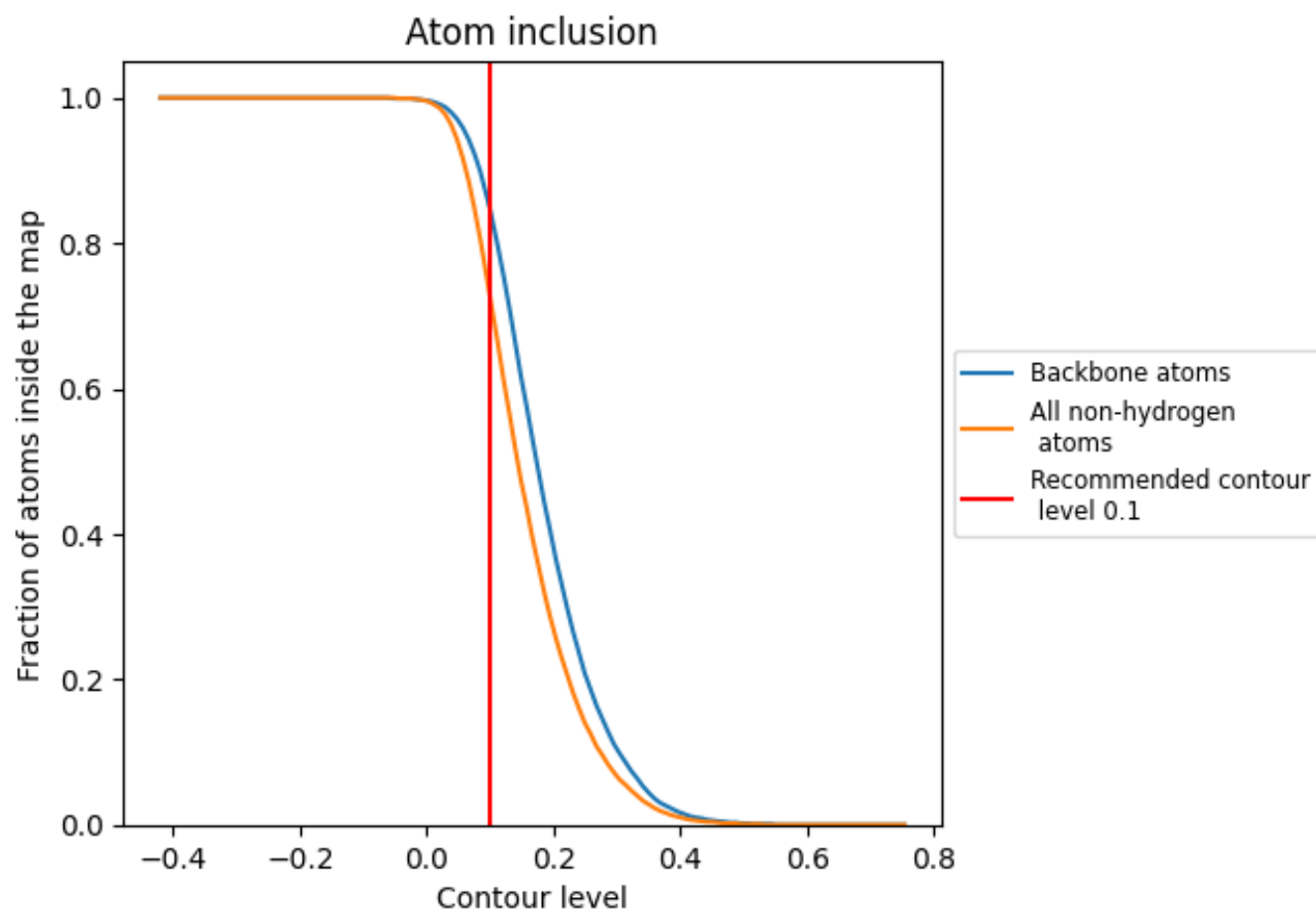
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 73% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7260	0.3580
A	0.7870	0.4110
B	0.7810	0.3810
C	0.7580	0.3740
D	0.6760	0.3300
E	0.7720	0.3720

