



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 07:45 PM UTC

PDB ID : 6EYD / pdb_00006eyd
EMDB ID : EMD-3983
Title : Structure of Mycobacterium smegmatis RNA polymerase Sigma-A holoenzyme
Authors : Kouba, T.; Barvik, I.; Krasny, L.
Deposited on : 2017-11-11
Resolution : 4.22 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

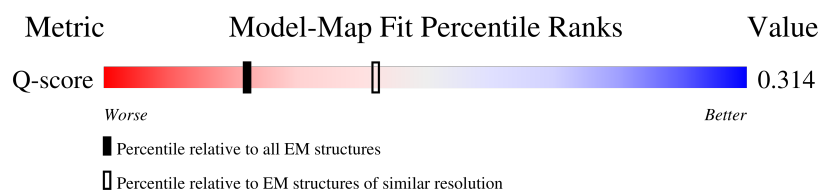
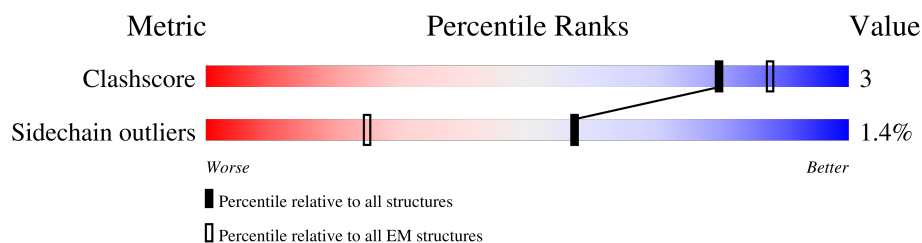
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.22 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	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	4772 (3.72 - 4.72)

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 22931 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	220	Total	C	N	O	S	0	0
			1617	1022	278	315	2		
1	B	233	Total	C	N	O	S	0	0
			1660	1050	287	321	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	1092	Total	C	N	O	S	0	0
			7982	5009	1395	1542	36		

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	1241	Total	C	N	O	S	0	0
			9399	5890	1702	1767	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 a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	217	Total	C	N	O	S	0	0
			1678	1060	300	312	6		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

SEQUENCE-PLOTS INFOmissingINFO

3 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	42579	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	0.537	Depositor
Minimum map value	-0.227	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.018	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

4 Model quality

4.1 Standard geometry

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	1.13	0/1643	1.46	0/2242
1	B	1.15	0/1686	1.51	3/2308 (0.1%)
2	C	1.11	0/8126	1.51	4/11071 (0.0%)
3	D	1.13	0/9549	1.57	4/12947 (0.0%)
4	E	1.07	0/604	1.59	0/822
5	F	1.14	0/1701	1.67	0/2301
All	All	1.12	0/23309	1.55	11/31691 (0.0%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	30	PHE	CA-CB-CG	5.56	119.36	113.80
2	C	135	THR	CA-C-N	5.29	125.12	120.10
2	C	135	THR	C-N-CA	5.29	125.12	120.10
3	D	915	VAL	CA-C-N	5.18	127.22	120.28
3	D	915	VAL	C-N-CA	5.18	127.22	120.28
3	D	505	HIS	CA-C-N	5.13	127.42	120.38
3	D	505	HIS	C-N-CA	5.13	127.42	120.38
1	B	212	GLY	CA-C-N	5.11	125.75	120.03
1	B	212	GLY	C-N-CA	5.11	125.75	120.03
2	C	886	ILE	CA-C-N	5.03	125.52	119.94
2	C	886	ILE	C-N-CA	5.03	125.52	119.94

There are no chirality outliers.

There are no planarity outliers.

4.2 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	1617	0	1616	14	0
1	B	1660	0	1624	14	0
2	C	7982	0	7577	50	0
3	D	9399	0	9220	49	0
4	E	592	0	583	3	0
5	F	1678	0	1671	6	0
6	D	2	0	0	0	0
7	D	1	0	0	0	0
All	All	22931	0	22291	120	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 (120) 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:B:26:LEU:HB3	1:B:30:PHE:HB3	1.60	0.82
2:C:882:ASN:HD21	2:C:1019:MET:HE1	1.57	0.70
1:A:48:GLY:HA2	1:A:170:PRO:HG3	1.77	0.66
1:B:110:ILE:HD11	1:B:118:VAL:HG21	1.83	0.60
2:C:678:ARG:O	2:C:695:ASP:N	2.33	0.60
2:C:602:MET:HE1	2:C:883:LYS:HG2	1.83	0.59
2:C:876:LEU:HD13	2:C:886:ILE:HD11	1.85	0.59
1:A:21:PHE:HB2	1:A:194:ILE:HG23	1.85	0.57
2:C:247:GLU:O	2:C:249:MET:N	2.37	0.57
3:D:883:VAL:HG11	3:D:1157:VAL:HG13	1.87	0.56
2:C:713:ALA:O	2:C:907:ILE:HA	2.06	0.55
5:F:173:ILE:HG22	5:F:238:LEU:HD13	1.90	0.54
1:B:15:ALA:HB3	1:B:18:ARG:HG2	1.90	0.53
3:D:589:THR:HG21	3:D:687:MET:HG2	1.89	0.53
2:C:739:THR:HG22	2:C:869:LYS:HA	1.91	0.53
2:C:1082:ILE:HG23	2:C:1106:PRO:HB3	1.91	0.53
3:D:407:LYS:HA	3:D:1230:THR:HG21	1.90	0.52
1:B:221:LEU:HD12	1:B:222:ALA:N	2.24	0.52
3:D:898:VAL:HG11	3:D:919:ALA:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:308:VAL:HA	5:F:311:VAL:HG12	1.92	0.52
1:A:9:LEU:HB3	1:B:221:LEU:HD13	1.92	0.52
2:C:696:GLY:N	2:C:699:THR:OG1	2.43	0.52
2:C:717:TRP:HA	3:D:724:THR:HG22	1.93	0.51
2:C:653:ALA:HB2	2:C:659:GLN:HE22	1.76	0.51
2:C:716:PRO:HA	2:C:721:ASN:HD21	1.75	0.51
2:C:913:VAL:HB	2:C:914:PRO:HD3	1.92	0.51
2:C:1028:VAL:O	2:C:1032:ILE:HG22	2.11	0.51
5:F:179:LEU:HD11	5:F:187:LEU:HD12	1.92	0.51
2:C:1093:VAL:HG23	2:C:1103:ILE:HG23	1.93	0.50
3:D:993:ALA:HB2	3:D:1261:ALA:HB1	1.93	0.50
2:C:1114:VAL:HG22	3:D:324:LEU:HD11	1.94	0.50
1:A:51:VAL:HG23	1:A:164:VAL:HG11	1.93	0.49
2:C:727:ILE:O	2:C:906:ILE:HA	2.12	0.49
3:D:567:SER:HB3	3:D:574:LEU:HD23	1.93	0.49
3:D:456:VAL:HA	3:D:490:VAL:HG11	1.94	0.49
1:A:95:MET:HE1	1:A:140:VAL:HG23	1.95	0.49
2:C:369:LEU:HD11	2:C:446:LEU:HD22	1.94	0.49
3:D:459:ARG:HG3	3:D:490:VAL:HG12	1.94	0.49
2:C:113:CYS:SG	2:C:120:TYR:N	2.86	0.49
1:B:22:VAL:HG12	1:B:193:ILE:HG22	1.93	0.48
3:D:675:LEU:HD22	3:D:714:LYS:HB3	1.96	0.48
2:C:30:VAL:HB	2:C:954:LEU:HD13	1.95	0.48
1:A:77:ILE:HG23	1:A:164:VAL:HG23	1.96	0.48
2:C:493:VAL:HG13	2:C:494:TYR:CD1	2.49	0.48
2:C:571:ASP:HA	2:C:576:GLN:HE22	1.78	0.48
1:A:85:VAL:HG21	1:A:140:VAL:HG21	1.96	0.47
1:B:43:LEU:HD11	3:D:620:MET:HG2	1.96	0.47
2:C:217:ILE:HD11	2:C:268:ILE:HD11	1.97	0.47
2:C:446:LEU:HD21	2:C:491:LEU:HD12	1.97	0.47
3:D:995:GLY:HA2	3:D:1157:VAL:HG11	1.97	0.47
2:C:45:LEU:HD21	2:C:155:GLY:HA3	1.97	0.46
2:C:887:GLY:HA2	3:D:431:VAL:HB	1.97	0.46
2:C:24:PRO:HB3	2:C:688:GLU:HB3	1.97	0.46
1:B:146:TYR:HB2	3:D:620:MET:HE2	1.98	0.46
2:C:1058:ARG:HA	3:D:421:ARG:HA	1.98	0.45
3:D:75:CYS:SG	3:D:77:ARG:HB2	2.57	0.45
3:D:663:ALA:HB1	3:D:665:THR:HG23	1.97	0.45
3:D:667:LEU:HD22	3:D:671:MET:HE2	1.98	0.45
3:D:965:LEU:HD13	3:D:1132:GLN:HB3	1.98	0.45
1:B:26:LEU:HB3	1:B:30:PHE:CB	2.38	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1065:TRP:CD1	3:D:877:VAL:HG11	2.51	0.45
3:D:320:ILE:HG21	3:D:340:LEU:HD23	1.97	0.45
3:D:486:VAL:O	3:D:490:VAL:HG13	2.17	0.45
3:D:1239:ILE:HG13	3:D:1240:ASN:HD22	1.82	0.45
3:D:556:ARG:HD3	4:E:89:LEU:HD13	1.99	0.45
5:F:239:ARG:HA	5:F:242:VAL:HG12	1.99	0.45
3:D:404:ASP:HA	3:D:407:LYS:HG2	1.99	0.44
4:E:41:LEU:HD21	4:E:47:LYS:HB3	1.99	0.44
1:A:95:MET:SD	1:A:95:MET:N	2.90	0.44
1:B:38:LEU:O	1:B:42:LEU:HD23	2.16	0.44
2:C:617:VAL:HG13	2:C:927:LEU:HD23	2.00	0.44
3:D:488:GLU:HA	3:D:491:ILE:HG22	1.99	0.44
1:A:78:LEU:HD21	2:C:611:ARG:HD3	2.00	0.43
2:C:1044:THR:HG21	2:C:1046:GLN:HE21	1.83	0.43
3:D:666:THR:HG22	3:D:669:ARG:HD2	2.00	0.43
3:D:366:ILE:HG23	5:F:260:GLN:HE22	1.83	0.43
2:C:794:VAL:HB	2:C:828:LEU:HB3	1.99	0.43
2:C:911:HIS:CE1	3:D:580:ASP:HA	2.53	0.43
2:C:45:LEU:HD11	2:C:443:LYS:HD2	2.01	0.43
3:D:277:LEU:HD11	3:D:295:ARG:HG2	2.00	0.43
3:D:1078:TYR:HB2	3:D:1081:LEU:HD21	2.00	0.43
2:C:468:PRO:HB2	2:C:557:GLY:HA3	2.00	0.43
2:C:575:ARG:O	2:C:575:ARG:HG3	2.19	0.43
1:B:41:THR:HG21	1:B:215:LEU:HG	2.01	0.42
3:D:1141:GLU:HA	3:D:1144:ARG:HD2	2.01	0.42
1:A:86:SER:O	1:A:116:VAL:HA	2.20	0.42
3:D:557:ILE:O	3:D:563:ASN:ND2	2.52	0.42
3:D:661:TRP:CZ3	3:D:663:ALA:HB2	2.55	0.42
3:D:876:LEU:HD23	3:D:1002:ILE:HG23	2.02	0.42
1:B:171:VAL:HA	1:B:198:THR:HA	2.01	0.42
2:C:166:VAL:HB	2:C:372:VAL:HG22	2.02	0.42
3:D:413:PHE:HA	3:D:417:LEU:HB2	2.01	0.42
3:D:896:ILE:HG21	3:D:965:LEU:HD12	2.01	0.42
3:D:1004:GLU:HG3	3:D:1005:PRO:HD3	2.00	0.42
1:A:39:ARG:HA	1:A:174:VAL:HG11	2.02	0.42
1:B:42:LEU:HD22	1:B:211:ALA:CB	2.50	0.42
2:C:1072:ALA:HA	3:D:554:GLU:CD	2.45	0.42
3:D:1029:LEU:HA	3:D:1032:VAL:HG22	2.02	0.41
1:A:48:GLY:CA	1:A:170:PRO:HG3	2.47	0.41
3:D:1259:ILE:O	3:D:1264:GLY:HA3	2.21	0.41
2:C:583:ALA:HB1	2:C:621:MET:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:677:GLN:HA	2:C:695:ASP:O	2.21	0.41
1:B:80:LEU:HD22	1:B:138:LEU:HD21	2.03	0.41
2:C:466:VAL:HG23	3:D:853:HIS:ND1	2.35	0.41
2:C:736:ASP:HA	2:C:739:THR:HG23	2.03	0.41
2:C:1062:MET:SD	2:C:1062:MET:N	2.92	0.41
3:D:1165:ARG:NH1	3:D:1209:MET:SD	2.94	0.41
2:C:1076:LEU:HD23	3:D:1259:ILE:HG13	2.03	0.41
3:D:1065:ILE:HD12	3:D:1112:LEU:HD13	2.03	0.41
1:A:62:GLU:O	1:A:73:VAL:HG13	2.21	0.40
2:C:132:ASN:HD21	2:C:400:VAL:HG13	1.86	0.40
2:C:732:LEU:HD22	2:C:737:VAL:HG21	2.03	0.40
2:C:928:GLY:HA3	2:C:1018:TYR:O	2.21	0.40
1:A:95:MET:CE	1:A:140:VAL:HG23	2.52	0.40
2:C:706:LEU:N	2:C:1020:TYR:OH	2.54	0.40
3:D:635:VAL:HG21	3:D:670:VAL:HG21	2.03	0.40
2:C:47:VAL:HG12	2:C:497:VAL:HG23	2.03	0.40
3:D:363:PRO:HD3	5:F:234:LEU:HD13	2.04	0.40
3:D:459:ARG:CG	3:D:490:VAL:HG12	2.51	0.40
4:E:43:ARG:NE	4:E:99:ASP:OD1	2.54	0.40

There are no symmetry-related clashes.

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

4.3.2 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	174/297 (59%)	173 (99%)	1 (1%)	78	80
1	B	153/297 (52%)	149 (97%)	4 (3%)	40	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	677/988 (68%)	669 (99%)	8 (1%)	63	73
3	D	951/426 (223%)	936 (98%)	15 (2%)	55	69
4	E	-	60 (97%)	2 (3%)	34	55
5	F	-	163 (99%)	1 (1%)	78	80
All	All	2181/2008 (109%)	2150 (99%)	31 (1%)	57	70

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	194	ILE
1	B	30	PHE
1	B	84	VAL
1	B	134	LEU
1	B	191	LYS
2	C	41	VAL
2	C	45	LEU
2	C	302	VAL
2	C	313	LEU
2	C	924	GLU
2	C	965	THR
2	C	1021	ILE
2	C	1024	LEU
3	D	260	SER
3	D	399	LEU
3	D	418	LEU
3	D	428	SER
3	D	490	VAL
3	D	574	LEU
3	D	619	ILE
3	D	667	LEU
3	D	731	SER
3	D	816	LEU
3	D	823	VAL
3	D	838	SER
3	D	876	LEU
3	D	957	THR
3	D	1209	MET
4	E	33	THR
4	E	41	LEU
5	F	351	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	151	GLN
1	B	129	ASN
2	C	132	ASN
2	C	366	ASN
2	C	410	ASN
2	C	659	GLN
2	C	677	GLN
2	C	882	ASN
2	C	1068	GLN
2	C	1125	ASN
3	D	145	HIS
3	D	146	ASN
3	D	165	GLN
3	D	368	ASN
3	D	464	ASN
3	D	888	HIS
3	D	932	ASN
3	D	1000	GLN
3	D	1134	HIS
3	D	1146	GLN
3	D	1191	ASN
3	D	1240	ASN
4	E	85	GLN
5	F	263	ASN

4.3.3 RNA [i](#)

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

4.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

4.6 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.

4.7 Other polymers

There are no such residues in this entry.

4.8 Polymer linkage issues

There are no chain breaks in this entry.

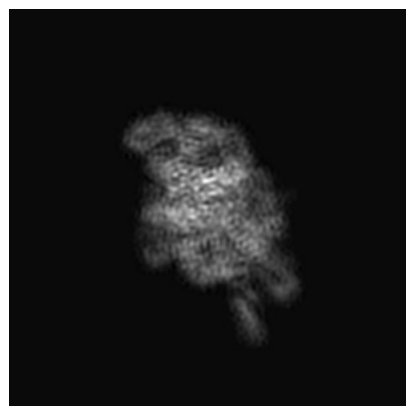
5 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3983. These allow visual inspection of the internal detail of the map and identification of artifacts.

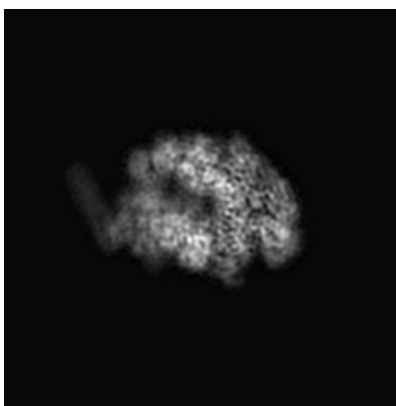
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

5.1 Orthogonal projections [i](#)

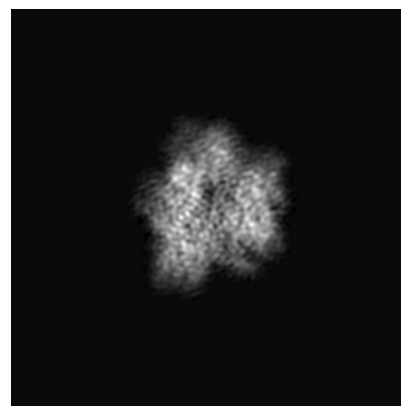
5.1.1 Primary map



X

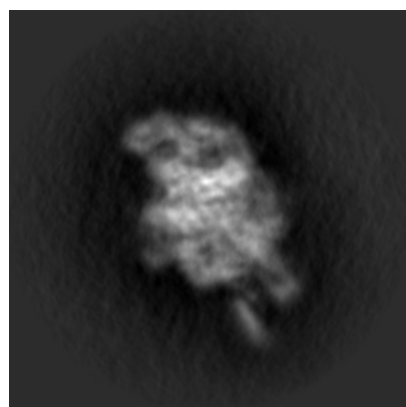


Y

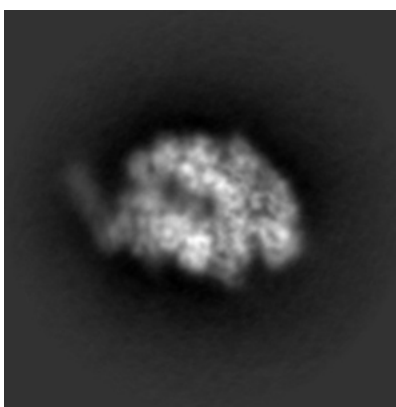


Z

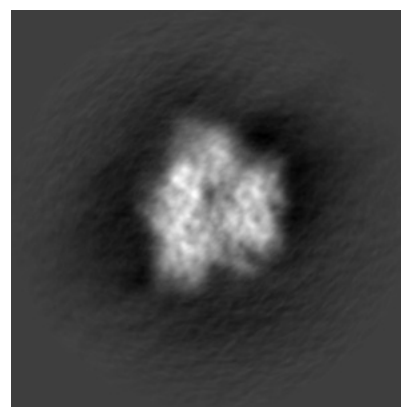
5.1.2 Raw map



X



Y

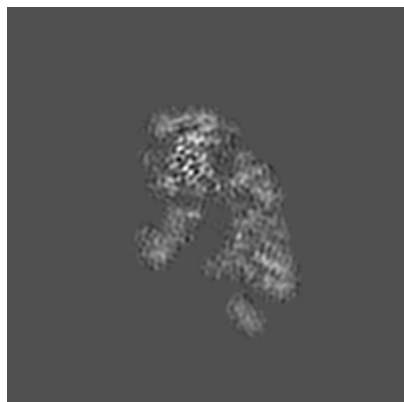


Z

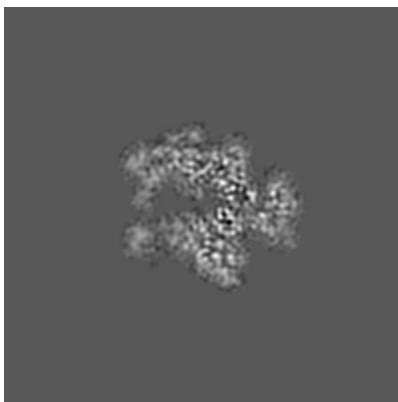
The images above show the map projected in three orthogonal directions.

5.2 Central slices [i](#)

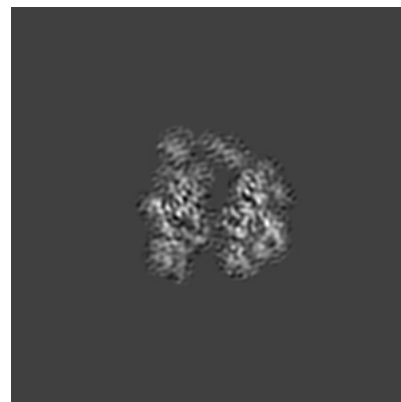
5.2.1 Primary map



X Index: 104

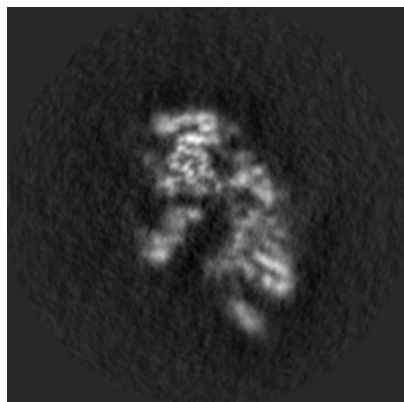


Y Index: 104

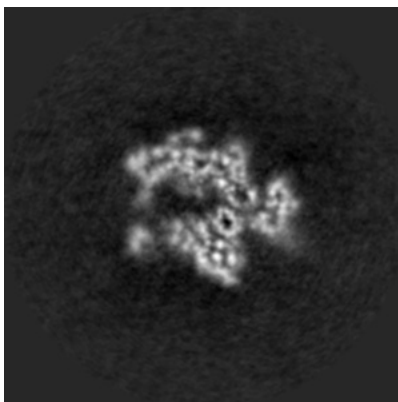


Z Index: 104

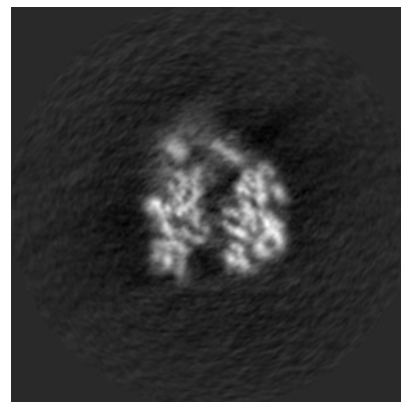
5.2.2 Raw map



X Index: 104



Y Index: 104

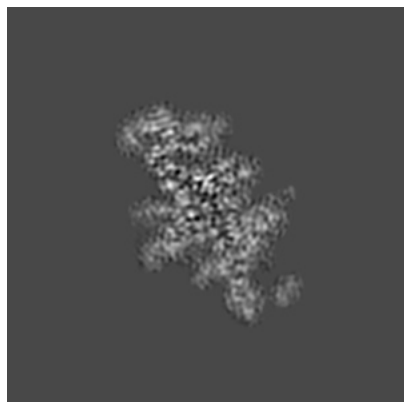


Z Index: 104

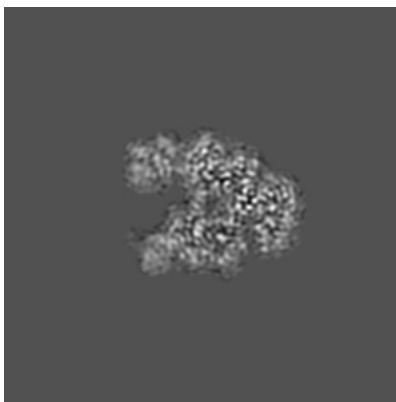
The images above show central slices of the map in three orthogonal directions.

5.3 Largest variance slices [i](#)

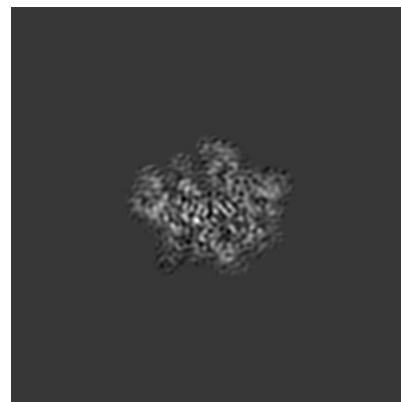
5.3.1 Primary map



X Index: 90

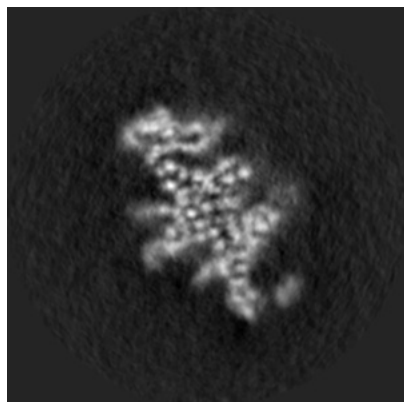


Y Index: 95

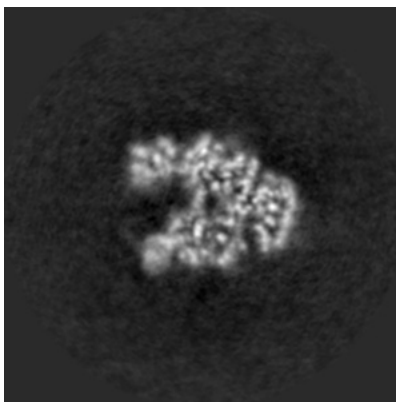


Z Index: 118

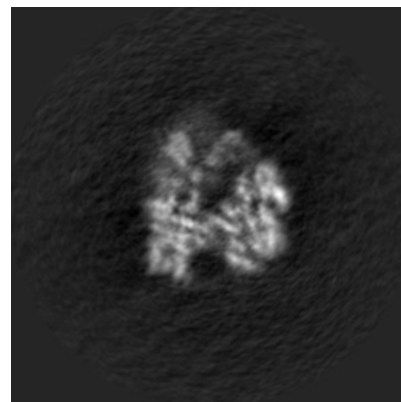
5.3.2 Raw map



X Index: 90



Y Index: 95

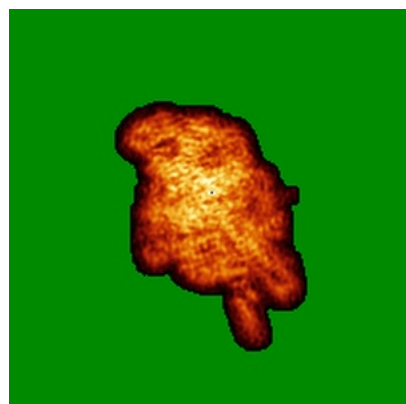


Z Index: 101

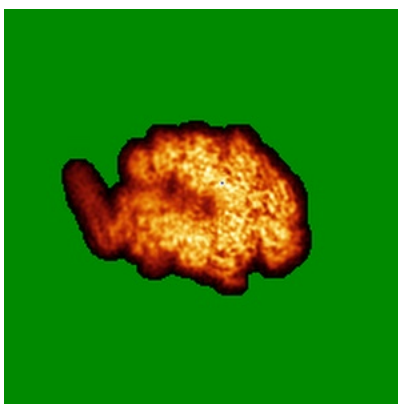
The images above show the largest variance slices of the map in three orthogonal directions.

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

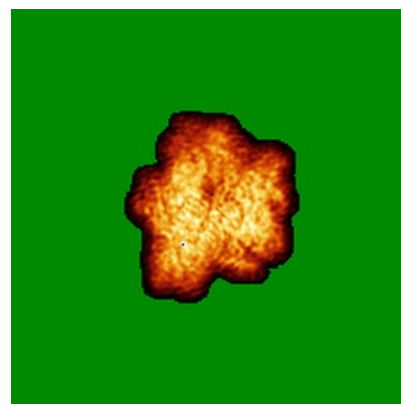
5.4.1 Primary map



X

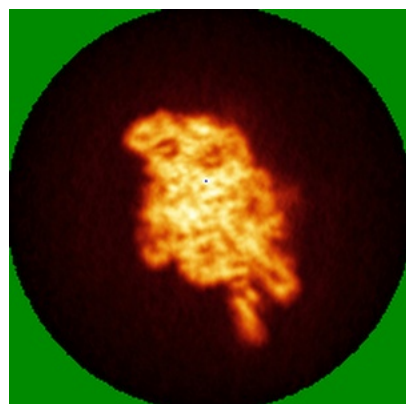


Y

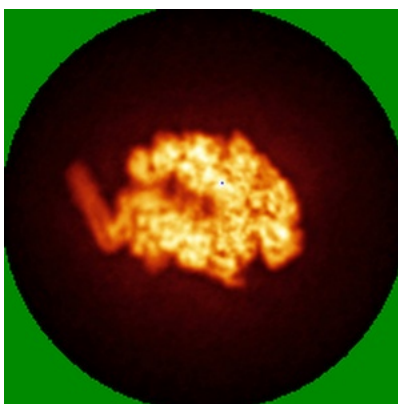


Z

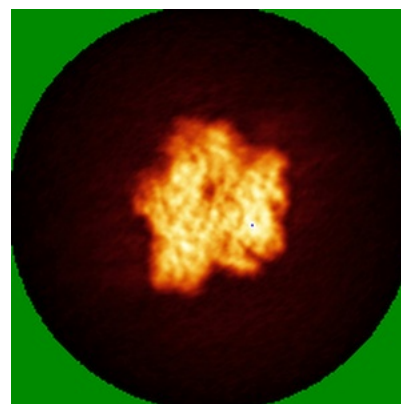
5.4.2 Raw map



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.

5.5 Orthogonal surface views [i](#)

5.5.1 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.

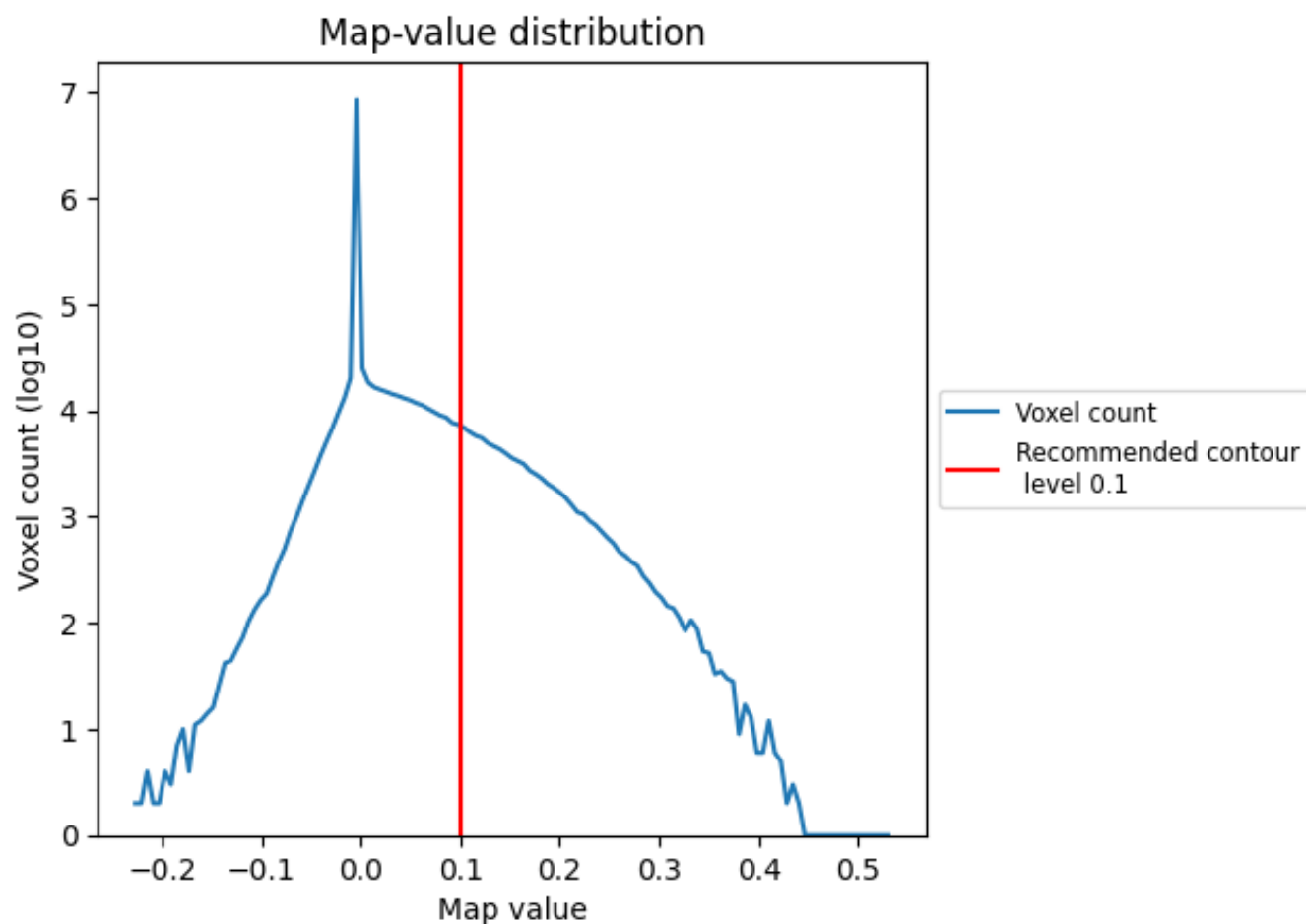
5.6 Mask visualisation [i](#)

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

6 Map analysis [i](#)

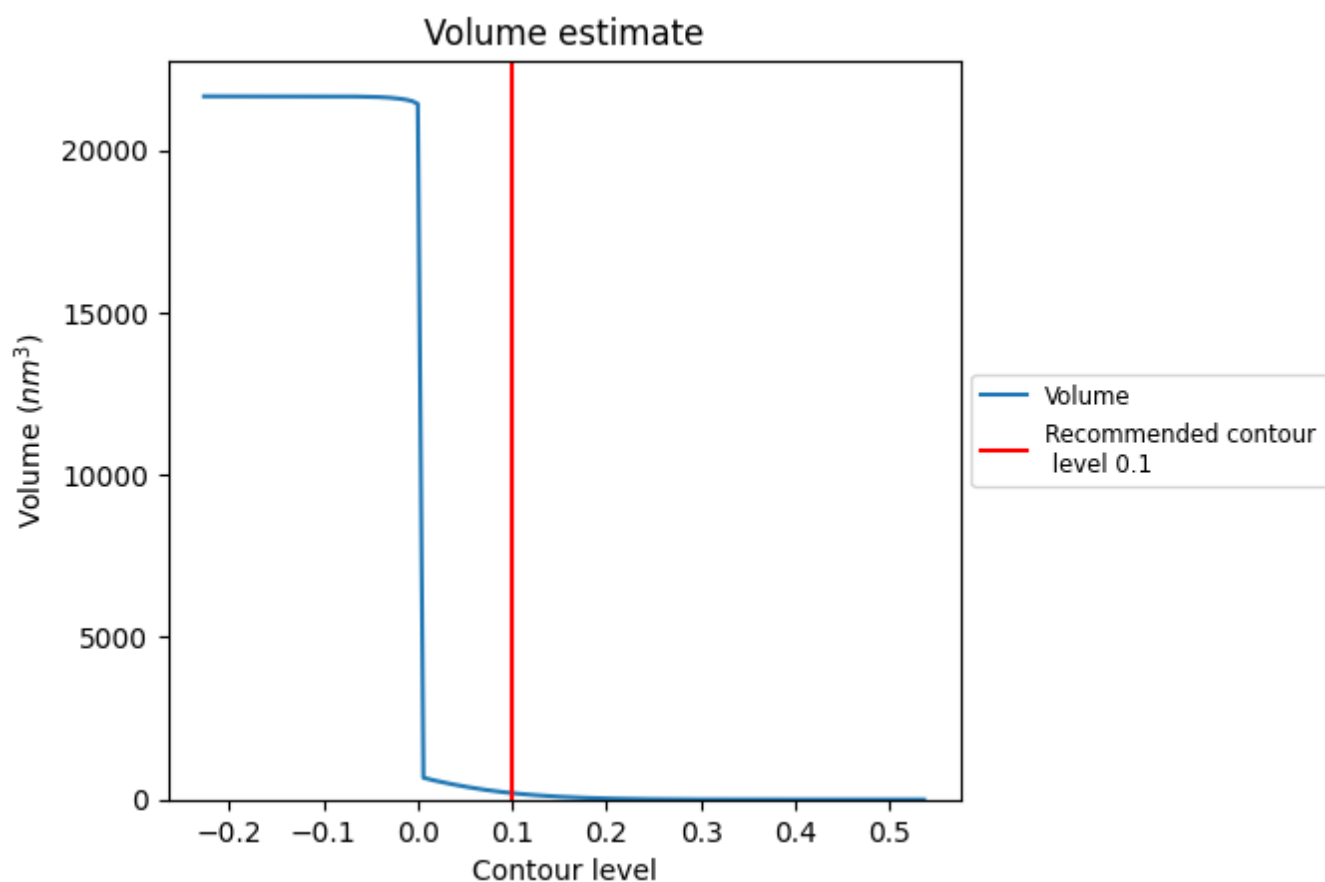
This section contains the results of statistical analysis of the map.

6.1 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.

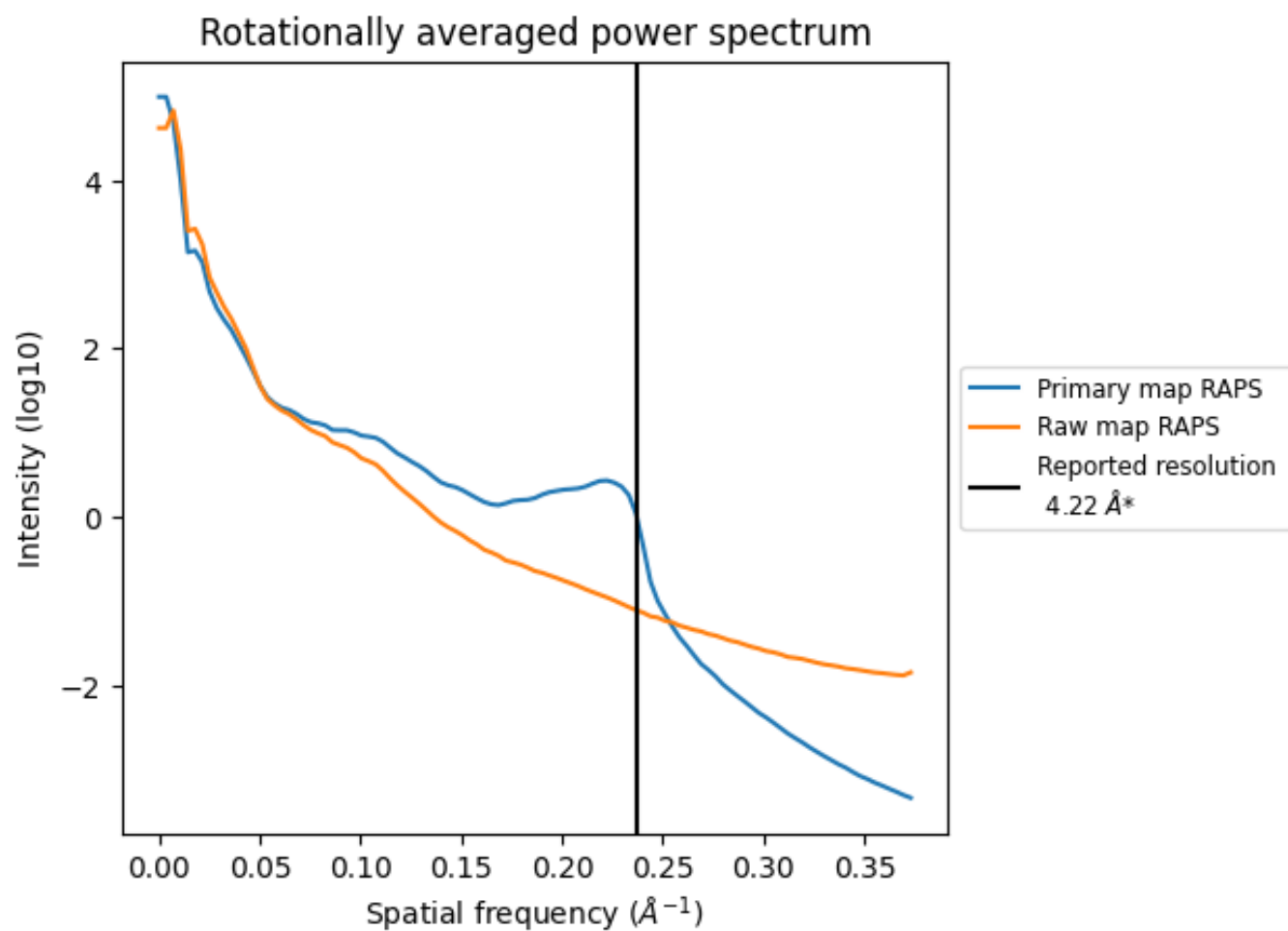
6.2 Volume estimate [i](#)



The volume at the recommended contour level is 199 nm³; this corresponds to an approximate mass of 180 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.

6.3 Rotationally averaged power spectrum ⓘ

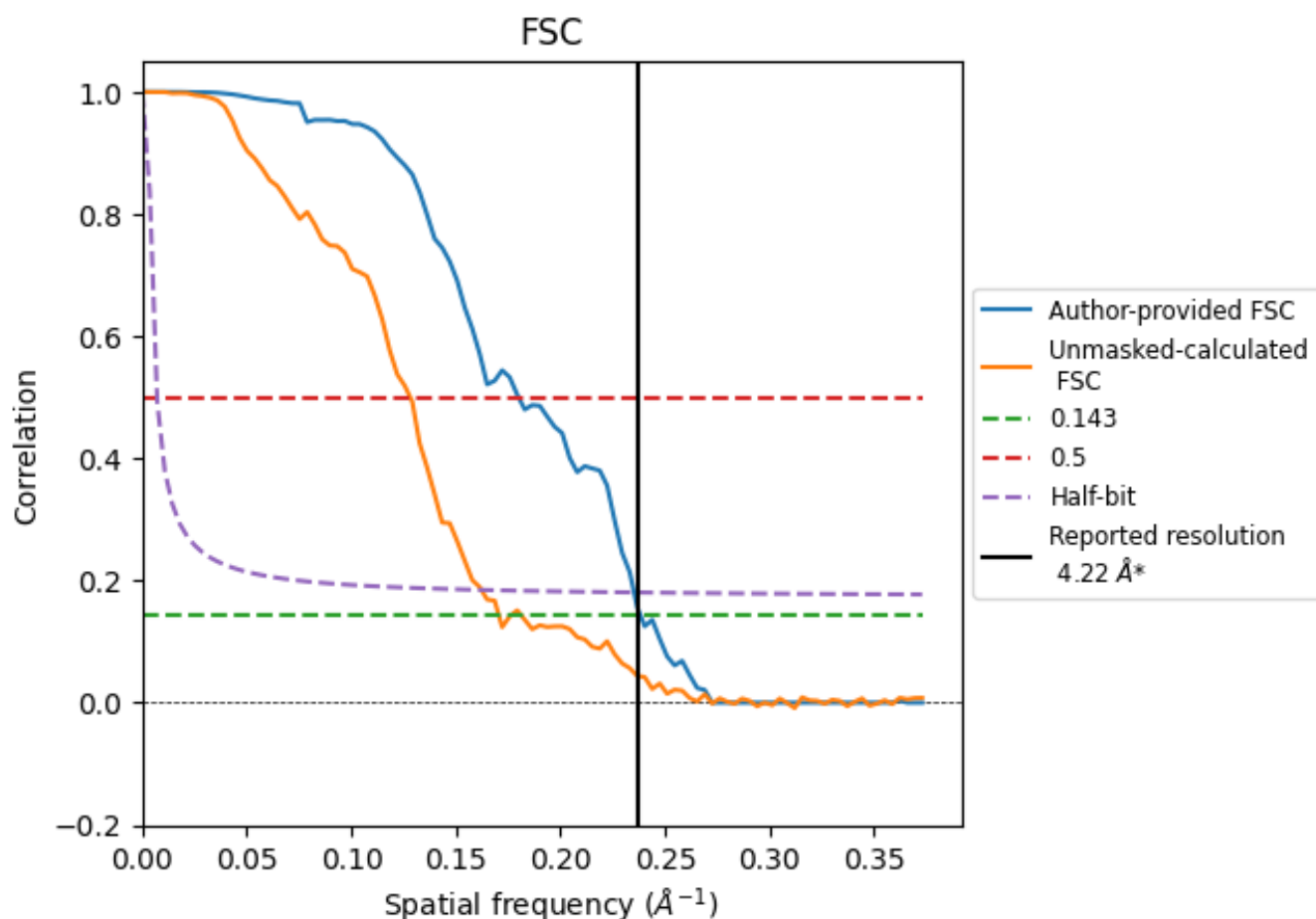


*Reported resolution corresponds to spatial frequency of 0.237 Å⁻¹

7 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.

7.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.237 \text{ \AA}^{-1}$

7.2 Resolution estimates [i](#)

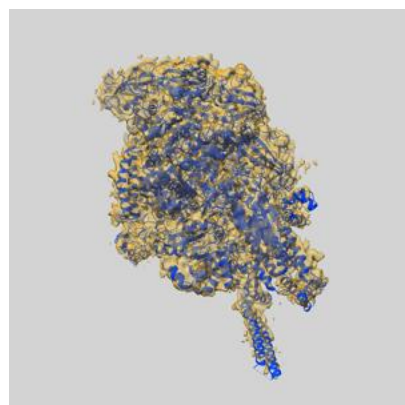
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.22	-	-
Author-provided FSC curve	4.20	5.56	4.25
Unmasked-calculated*	5.86	7.81	6.17

*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 5.86 differs from the reported value 4.22 by more than 10 %

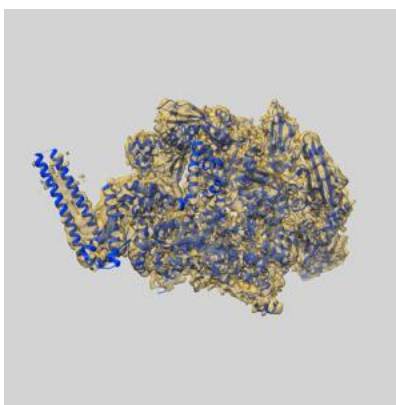
8 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-3983 and PDB model 6EYD. Per-residue inclusion information can be found in section ?? on page ??.

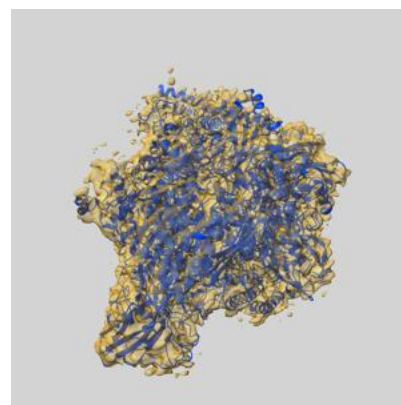
8.1 Map-model overlay [i](#)



X



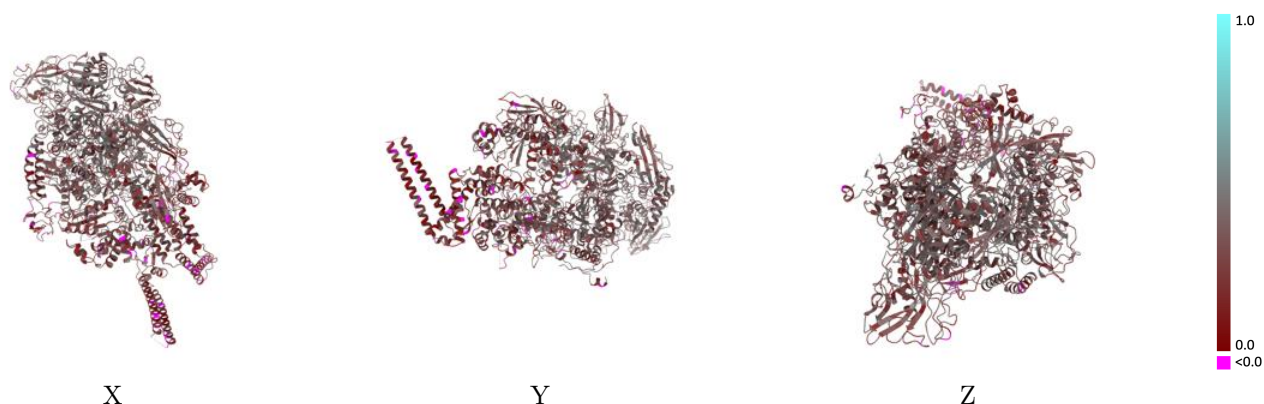
Y



Z

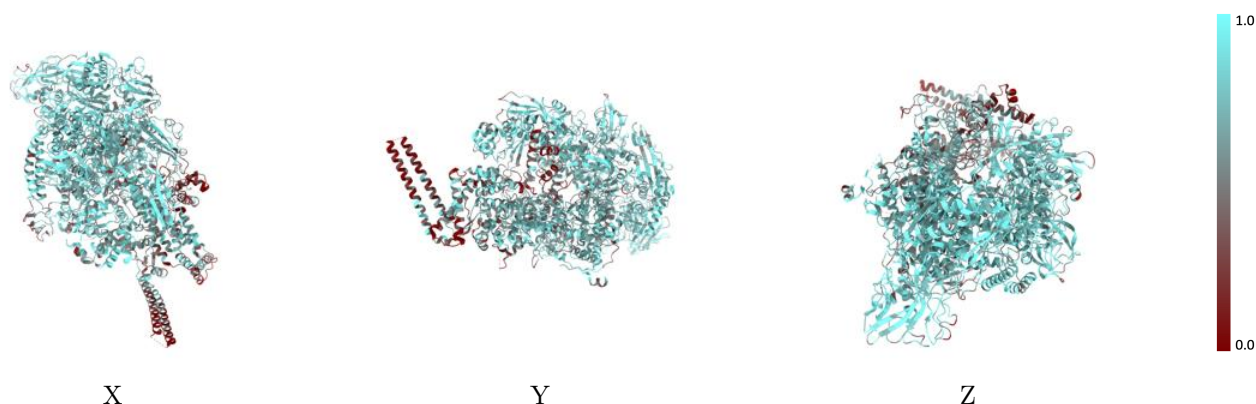
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



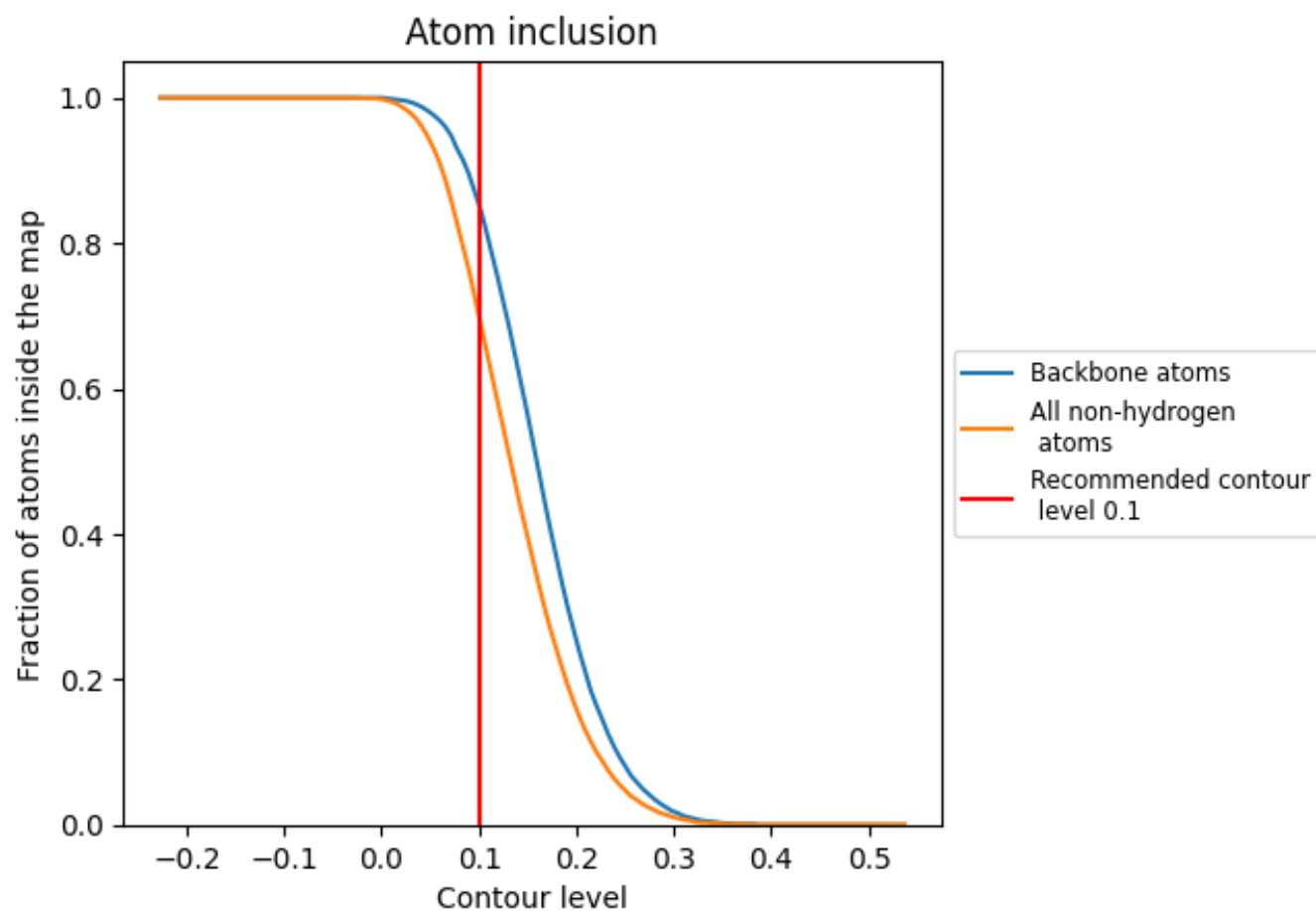
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

8.3 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).

8.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 70% of all non-hydrogen atoms, are inside the map.

8.5 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	0.7000	0.3140
A	0.7580	0.3570
B	0.7930	0.3420
C	0.7460	0.3350
D	0.6800	0.2980
E	0.7350	0.3230
F	0.4340	0.2260

1.0

0.0

<0.0