



Full wwPDB EM Validation Report ⓘ

Nov 4, 2024 – 02:30 AM JST

PDB ID : 7CKM
EMDB ID : EMD-30388
Title : Structure of Machupo virus polymerase bound to Z matrix protein (monomeric complex)
Authors : Xu, X.; Peng, R.; Peng, Q.; Shi, Y.
Deposited on : 2020-07-17
Resolution : 3.37 Å (reported)
Based on initial models : 5I72, 6KLD

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.dev113
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

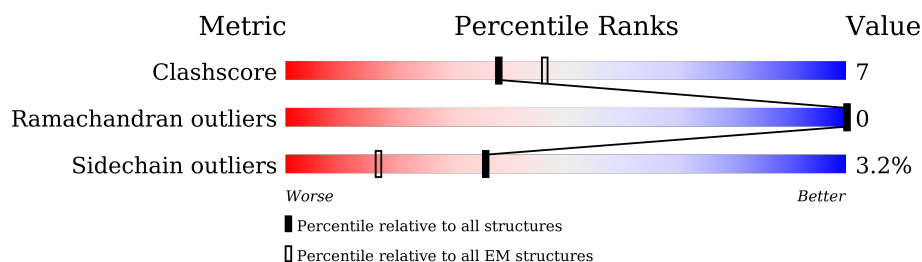
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.37 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	2209	
2	B	94	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13114 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 RNA-directed RNA polymerase L.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1603	Total	C	N	O	S	0	0
			12702	8101	2132	2384	85		

- Molecule 2 is a protein called RING finger protein Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	49	Total	C	N	O	S	0	0
			407	257	73	68	9		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	Mn	0
			1	1	

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	A	2	Total	Zn	0
			2	2	
4	B	2	Total	Zn	0
			2	2	



T81	ILE	THR	VAL	PRO	VAL	GLU	PRO	SER	ALA	PRO	PRO	PRO	MET	GLY	ASN	CYS	ASN	LYS	PRO	PRO	LYS	ARG	PRO	PRO	ASN	THR	GLN	THR	SER	SER	GLN	ASN	PRO	SER	ALA	GLU	PHE	ARG	ARG	THR	ALA	PRO	PRO	SER	L33	C39	A45	D46	T47	N48	C52	H55	Y56	L57	C58	L59	R60	T64	M65	L66	L71	C72	H73	I74	C75
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	370690	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.161	Depositor
Minimum map value	-0.103	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	259.2, 259.2, 259.2	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	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, MN

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	0.26	0/12922	0.43	0/17432
2	B	0.37	0/419	0.57	0/569
All	All	0.26	0/13341	0.43	0/18001

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	12702	0	12716	174	0
2	B	407	0	383	15	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
All	All	13114	0	13099	188	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 (188) 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:B:39:CYS:HB2	2:B:58:CYS:SG	1.96	1.06
2:B:45:ALA:HB3	2:B:58:CYS:SG	2.00	1.01
1:A:236:ARG:HG2	1:A:1316:LEU:HD11	1.71	0.73
1:A:456:PRO:HG2	1:A:525:LEU:HB2	1.71	0.73
2:B:45:ALA:CB	2:B:58:CYS:SG	2.77	0.72
1:A:274:GLY:HA3	1:A:672:VAL:HG21	1.72	0.72
1:A:1730:GLN:HG2	1:A:1731:LEU:H	1.58	0.68
1:A:60:CYS:SG	1:A:61:GLN:N	2.67	0.67
1:A:788:ASN:ND2	1:A:1209:GLU:OE1	2.30	0.65
2:B:55:HIS:HE1	2:B:75:CYS:HB3	1.60	0.65
1:A:204:LEU:HB2	1:A:207:LEU:HD23	1.79	0.65
1:A:412:ASN:O	1:A:416:ASN:ND2	2.30	0.65
1:A:55:CYS:O	1:A:60:CYS:HB3	1.97	0.64
1:A:1752:LEU:HG	1:A:1753:THR:HG23	1.79	0.62
2:B:59:LEU:O	2:B:59:LEU:HD12	1.99	0.62
1:A:694:LYS:O	1:A:698:ASN:ND2	2.32	0.62
1:A:285:PRO:HG2	1:A:615:LEU:HD21	1.81	0.62
1:A:539:LEU:HA	1:A:560:SER:O	2.00	0.62
1:A:1650:CYS:SG	1:A:1651:THR:N	2.72	0.62
1:A:11:LEU:HD23	1:A:163:LEU:HD11	1.82	0.62
1:A:638:PHE:HB2	1:A:655:LEU:HD21	1.82	0.62
1:A:1465:VAL:HG22	1:A:1466:LYS:H	1.65	0.61
1:A:1425:GLN:O	1:A:1429:ASN:ND2	2.34	0.61
2:B:55:HIS:CE1	2:B:75:CYS:HB3	2.35	0.61
1:A:382:VAL:HG23	1:A:385:MET:HB2	1.81	0.61
1:A:340:LEU:HB3	1:A:342:VAL:HG13	1.83	0.60
1:A:737:ASN:ND2	1:A:1282:SER:OG	2.35	0.60
1:A:1211:ARG:HD3	1:A:1216:ARG:HD2	1.84	0.59
1:A:186:ARG:NH2	1:A:1099:GLU:O	2.34	0.59
1:A:775:VAL:HG23	1:A:1365:LEU:HA	1.83	0.59
1:A:1388:VAL:HG12	1:A:1389:MET:H	1.66	0.59
1:A:1630:VAL:HG23	1:A:1638:LEU:HG	1.85	0.59
1:A:1419:ILE:HG21	1:A:1443:VAL:HG21	1.85	0.58
1:A:549:GLY:O	1:A:553:ARG:NH2	2.37	0.58
1:A:687:LEU:HD21	2:B:64:THR:HG21	1.87	0.57
1:A:1692:GLU:HB2	1:A:1788:ILE:HD11	1.86	0.57
1:A:1650:CYS:SG	1:A:1667:SER:OG	2.63	0.56
1:A:124:LEU:HD21	1:A:838:GLU:HG2	1.87	0.56
1:A:1138:THR:O	1:A:1142:THR:HG23	2.06	0.56
1:A:1402:ALA:O	1:A:1406:ASN:ND2	2.39	0.56
1:A:330:VAL:HG21	1:A:360:LEU:HD13	1.88	0.55
1:A:553:ARG:NH1	1:A:555:TYR:OH	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:VAL:HG12	1:A:737:ASN:H	1.72	0.54
1:A:60:CYS:HB2	1:A:101:LEU:HD21	1.89	0.54
1:A:1182:ASP:OD2	1:A:1336:LYS:NZ	2.39	0.54
2:B:55:HIS:HE1	2:B:75:CYS:CB	2.18	0.54
1:A:69:VAL:HA	1:A:72:ILE:HG22	1.89	0.54
1:A:718:ILE:O	1:A:722:GLU:HG2	2.08	0.53
1:A:1741:ARG:HH12	1:A:1763:ARG:HA	1.73	0.53
1:A:156:VAL:O	1:A:160:ILE:HG12	2.09	0.52
1:A:1465:VAL:HG21	1:A:1811:ALA:HA	1.90	0.52
1:A:1555:PHE:HD2	1:A:1632:ASP:HA	1.75	0.52
1:A:159:ARG:O	1:A:163:LEU:HG	2.09	0.52
1:A:539:LEU:HB3	1:A:561:ASP:HA	1.91	0.52
1:A:334:CYS:HA	1:A:337:ILE:HG22	1.92	0.51
1:A:93:LEU:HD12	1:A:98:LEU:HD13	1.91	0.51
1:A:898:THR:HG22	1:A:899:GLU:H	1.76	0.51
2:B:52:CYS:HB3	2:B:55:HIS:NE2	2.25	0.51
1:A:34:HIS:CD2	1:A:35:PRO:HD2	2.46	0.51
1:A:34:HIS:HB3	1:A:37:PRO:HD2	1.93	0.51
1:A:83:LEU:HD11	1:A:130:LEU:HD13	1.92	0.51
1:A:543:LEU:HD12	1:A:557:VAL:HG22	1.93	0.51
1:A:1093:ASN:HA	1:A:1097:TYR:HB3	1.92	0.51
1:A:737:ASN:HD21	1:A:1282:SER:H	1.59	0.51
1:A:1377:THR:O	1:A:1377:THR:OG1	2.28	0.50
1:A:908:ILE:O	1:A:912:VAL:HG23	2.12	0.50
1:A:971:ASP:N	1:A:971:ASP:OD1	2.44	0.50
1:A:864:SER:O	1:A:867:TYR:HB3	2.12	0.50
1:A:1195:THR:HG21	1:A:1368:PHE:HE2	1.77	0.50
1:A:1661:ILE:HG22	1:A:1662:ILE:HG12	1.94	0.49
1:A:42:GLY:O	1:A:45:LEU:HG	2.12	0.49
1:A:910:HIS:O	1:A:914:VAL:HG23	2.12	0.49
1:A:920:ALA:HB2	1:A:1007:LEU:HG	1.95	0.49
1:A:152:VAL:HG12	1:A:153:ALA:H	1.77	0.49
1:A:860:LEU:HA	1:A:1006:PRO:HA	1.95	0.49
1:A:982:GLU:N	1:A:982:GLU:OE1	2.43	0.49
1:A:664:GLU:OE1	1:A:664:GLU:N	2.46	0.49
1:A:1177:ALA:HB2	1:A:1334:LEU:HD23	1.95	0.49
1:A:165:ILE:HG23	1:A:191:LEU:HD23	1.95	0.49
1:A:681:GLY:HA3	1:A:685:ALA:HB2	1.93	0.49
1:A:261:THR:HG23	1:A:262:ASP:H	1.77	0.48
1:A:525:LEU:HD22	1:A:528:CYS:HB3	1.94	0.48
1:A:623:SER:O	1:A:623:SER:OG	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1243:ALA:HB1	1:A:1280:ILE:HD11	1.95	0.48
1:A:29:PHE:HD2	1:A:30:LEU:HD22	1.79	0.48
1:A:1109:SER:OG	1:A:1110:SER:N	2.46	0.47
1:A:580:LEU:O	1:A:584:SER:OG	2.30	0.47
1:A:279:SER:OG	1:A:669:ARG:NH1	2.47	0.47
1:A:366:ASP:OD1	1:A:367:SER:N	2.47	0.47
1:A:898:THR:O	1:A:902:VAL:HG23	2.14	0.47
2:B:39:CYS:HA	2:B:56:TYR:O	2.14	0.47
1:A:1654:ILE:HD13	1:A:1654:ILE:HA	1.82	0.47
2:B:46:ASP:N	2:B:46:ASP:OD1	2.48	0.47
1:A:508:ARG:HE	1:A:509:LEU:H	1.62	0.47
1:A:1733:MET:HG3	1:A:1734:ASN:H	1.80	0.47
1:A:215:ASP:N	1:A:215:ASP:OD1	2.47	0.47
1:A:969:ASP:N	1:A:969:ASP:OD1	2.44	0.47
1:A:1741:ARG:NH1	1:A:1763:ARG:HA	2.30	0.46
1:A:191:LEU:HD12	1:A:191:LEU:HA	1.82	0.46
2:B:47:THR:HG22	2:B:48:ASN:ND2	2.31	0.46
1:A:1235:VAL:HG13	1:A:1283:HIS:HB3	1.98	0.46
1:A:8:LEU:O	1:A:12:ILE:HG13	2.16	0.46
1:A:350:LEU:HD22	1:A:392:ARG:HH12	1.81	0.46
1:A:1291:GLY:HA3	1:A:1294:ILE:HD12	1.96	0.46
1:A:457:ILE:HD12	1:A:459:ARG:NE	2.31	0.45
1:A:1110:SER:O	1:A:1110:SER:OG	2.31	0.45
1:A:1351:TRP:O	1:A:1355:ILE:HG12	2.16	0.45
1:A:233:SER:O	1:A:237:GLU:HG2	2.15	0.45
1:A:588:LEU:HA	1:A:591:MET:HG2	1.98	0.45
1:A:68:PHE:CE1	1:A:85:LYS:HB3	2.52	0.45
1:A:498:ILE:HG22	1:A:1619:ILE:HD11	1.98	0.45
1:A:321:SER:O	1:A:321:SER:OG	2.30	0.45
1:A:459:ARG:HD3	1:A:522:ASN:ND2	2.32	0.45
1:A:1364:LYS:HD3	1:A:1364:LYS:HA	1.68	0.45
1:A:230:ILE:N	1:A:785:SER:OG	2.48	0.45
1:A:1161:LEU:HD11	1:A:1296:HIS:CE1	2.52	0.45
1:A:1149:SER:OG	1:A:1298:THR:HG22	2.18	0.44
1:A:798:LEU:H	1:A:798:LEU:HD23	1.82	0.44
1:A:226:LEU:HD21	1:A:1203:ALA:HB1	1.99	0.44
1:A:502:LYS:HG3	1:A:1619:ILE:HG23	1.99	0.44
1:A:547:LYS:HB2	1:A:553:ARG:HD3	2.00	0.44
1:A:722:GLU:OE2	1:A:1266:SER:HB3	2.17	0.44
1:A:117:LYS:NZ	1:A:1135:ASP:OD1	2.48	0.44
1:A:97:THR:HG21	1:A:919:MET:HG3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:LYS:HD2	1:A:649:ILE:HG13	1.98	0.44
1:A:1200:LEU:HD21	1:A:1302:TYR:CZ	2.53	0.44
1:A:443:SER:O	1:A:447:ILE:HG13	2.18	0.44
1:A:545:TYR:HB3	1:A:555:TYR:CD1	2.52	0.44
1:A:1223:SER:O	1:A:1227:ILE:HG12	2.18	0.44
1:A:1730:GLN:HG2	1:A:1731:LEU:N	2.28	0.44
1:A:96:LYS:HD2	1:A:135:VAL:HG12	1.99	0.44
1:A:261:THR:HA	1:A:647:HIS:CE1	2.53	0.44
1:A:786:LEU:HA	1:A:786:LEU:HD12	1.81	0.44
2:B:57:LEU:HD12	2:B:57:LEU:HA	1.76	0.44
1:A:840:LEU:HD11	1:A:872:ASN:HB2	1.99	0.43
1:A:1286:SER:OG	1:A:1287:VAL:N	2.50	0.43
1:A:1237:GLU:OE2	1:A:1283:HIS:NE2	2.51	0.43
1:A:870:LEU:HB3	1:A:893:MET:HE3	2.01	0.43
1:A:440:ASP:OD1	1:A:440:ASP:N	2.39	0.43
1:A:34:HIS:O	1:A:38:LEU:HB2	2.19	0.43
1:A:766:ASP:N	1:A:766:ASP:OD2	2.52	0.43
1:A:1783:ASN:HD22	1:A:1786:LEU:HD12	1.83	0.43
1:A:218:LEU:O	1:A:222:LEU:HB2	2.19	0.42
1:A:354:ASP:O	1:A:358:VAL:HG22	2.19	0.42
1:A:1326:THR:HG22	1:A:1328:SER:H	1.83	0.42
1:A:1688:THR:HG22	1:A:1788:ILE:HD12	2.01	0.42
1:A:1552:LEU:HD12	1:A:1630:VAL:HG11	2.01	0.42
1:A:1774:ASP:OD2	1:A:1775:VAL:N	2.52	0.42
1:A:832:LEU:HD13	1:A:1566:LEU:HD23	2.00	0.42
1:A:83:LEU:HB3	1:A:126:LEU:HD11	2.02	0.42
1:A:36:SER:HB3	1:A:37:PRO:HD3	2.00	0.42
1:A:335:ASN:HA	1:A:389:VAL:HG23	2.02	0.42
1:A:1167:PHE:CE1	1:A:1383:LYS:HD2	2.55	0.42
1:A:1699:SER:OG	1:A:1700:THR:N	2.52	0.42
1:A:1209:GLU:HB2	1:A:1218:LYS:HG3	2.02	0.42
1:A:1465:VAL:HG22	1:A:1466:LYS:N	2.33	0.42
1:A:1469:VAL:HG12	1:A:1801:TYR:CG	2.55	0.42
1:A:261:THR:HG23	1:A:262:ASP:N	2.35	0.41
1:A:745:THR:O	1:A:749:GLU:HG2	2.20	0.41
1:A:1173:ASP:OD2	1:A:1325:TYR:OH	2.28	0.41
1:A:19:ARG:HG3	1:A:52:LEU:HD11	2.03	0.41
2:B:71:LEU:HD12	2:B:71:LEU:HA	1.94	0.41
1:A:457:ILE:HD12	1:A:459:ARG:HE	1.85	0.41
1:A:1142:THR:HG22	1:A:1231:HIS:CE1	2.55	0.41
1:A:101:LEU:HA	1:A:140:ILE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1110:SER:HB2	1:A:1230:TRP:CH2	2.55	0.41
1:A:1115:LEU:HD22	1:A:1138:THR:HB	2.03	0.41
1:A:114:LYS:HE2	1:A:114:LYS:HB2	1.92	0.41
1:A:1211:ARG:HG3	1:A:1216:ARG:HA	2.02	0.41
1:A:4:TYR:HB2	1:A:167:ILE:HD11	2.02	0.41
1:A:29:PHE:HD1	1:A:41:GLU:HB3	1.86	0.41
1:A:521:GLY:O	1:A:522:ASN:ND2	2.54	0.41
1:A:687:LEU:HD12	1:A:687:LEU:HA	1.89	0.41
1:A:1690:ALA:HA	1:A:1694:GLY:O	2.20	0.41
1:A:8:LEU:HG	1:A:12:ILE:HD11	2.01	0.41
1:A:96:LYS:O	1:A:135:VAL:HB	2.21	0.41
1:A:372:ILE:HD11	1:A:447:ILE:HG23	2.03	0.41
1:A:786:LEU:HD21	1:A:1312:TYR:CE2	2.56	0.41
1:A:1531:LEU:HB3	1:A:1536:CYS:SG	2.61	0.40
1:A:1588:LYS:HE3	1:A:1588:LYS:HB3	1.96	0.40
1:A:48:SER:HA	1:A:51:GLU:HG3	2.03	0.40
1:A:273:LEU:HB2	1:A:656:VAL:HG12	2.03	0.40
1:A:9:LYS:HB3	1:A:30:LEU:HD13	2.04	0.40
1:A:1272:GLN:O	1:A:1276:LEU:HB2	2.22	0.40
1:A:1753:THR:N	1:A:1754:PRO:HD3	2.37	0.40
1:A:19:ARG:HB3	1:A:21:GLU:OE1	2.22	0.40
1:A:234:THR:O	1:A:238:GLU:HG2	2.21	0.40
2:B:72:CYS:SG	2:B:74:ILE:HG22	2.61	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	1569/2209 (71%)	1441 (92%)	128 (8%)	0	100	100
2	B	47/94 (50%)	36 (77%)	11 (23%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1616/2303 (70%)	1477 (91%)	139 (9%)	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	1443/2007 (72%)	1400 (97%)	43 (3%)	36	62
2	B	47/88 (53%)	42 (89%)	5 (11%)	5	20
All	All	1490/2095 (71%)	1442 (97%)	48 (3%)	36	60

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	PHE
1	A	60	CYS
1	A	108	ASN
1	A	110	ASP
1	A	155	TRP
1	A	201	LEU
1	A	204	LEU
1	A	234	THR
1	A	382	VAL
1	A	406	LYS
1	A	421	LEU
1	A	451	LEU
1	A	564	VAL
1	A	585	ASP
1	A	691	ARG
1	A	772	THR
1	A	798	LEU
1	A	865	THR
1	A	980	TYR
1	A	1010	LEU

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Mol	Chain	Res	Type
1	A	1125	VAL
1	A	1138	THR
1	A	1162	ASN
1	A	1212	THR
1	A	1288	LEU
1	A	1290	MET
1	A	1296	HIS
1	A	1330	ASP
1	A	1352	LEU
1	A	1366	ASN
1	A	1377	THR
1	A	1397	THR
1	A	1426	CYS
1	A	1466	LYS
1	A	1559	ARG
1	A	1560	LEU
1	A	1638	LEU
1	A	1654	ILE
1	A	1688	THR
1	A	1717	TRP
1	A	1731	LEU
1	A	1755	PHE
1	A	1795	THR
2	B	39	CYS
2	B	57	LEU
2	B	59	LEU
2	B	60	ARG
2	B	66	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	384	GLN
1	A	522	ASN
1	A	628	ASN
1	A	647	HIS
1	A	737	ASN
1	A	788	ASN
1	A	1013	ASN
1	A	1128	ASN
1	A	1206	GLN

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Mol	Chain	Res	Type
1	A	1296	HIS
1	A	1308	GLN
1	A	1366	ASN
1	A	1429	ASN
1	A	1542	ASN
1	A	1611	GLN
1	A	1783	ASN
2	B	48	ASN

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](#)

Of 5 ligands modelled in this entry, 5 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

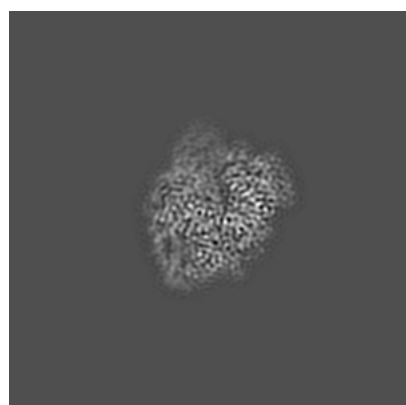
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30388. These allow visual inspection of the internal detail of the map and identification of artifacts.

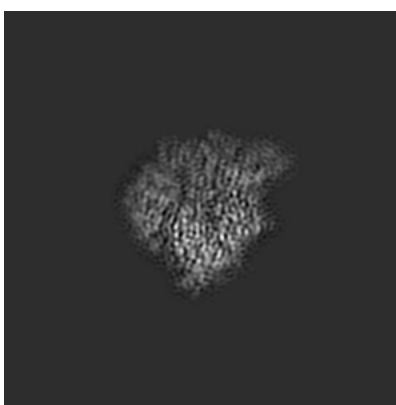
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

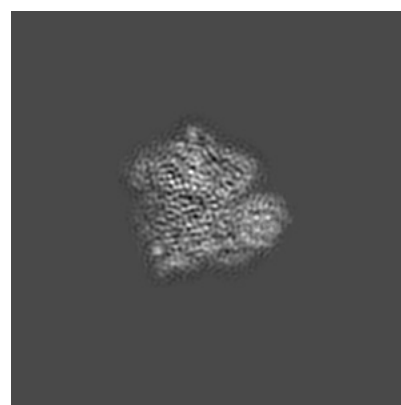
6.1.1 Primary map



X



Y

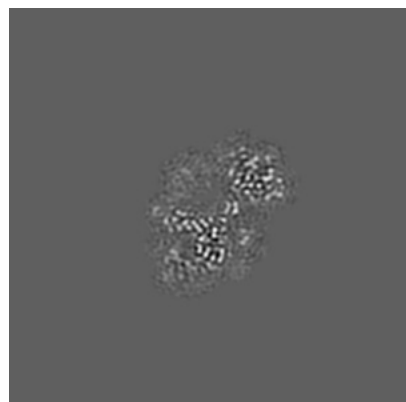


Z

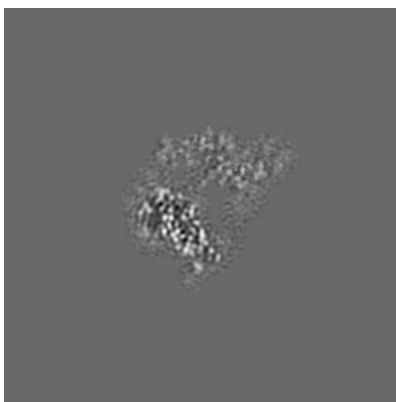
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

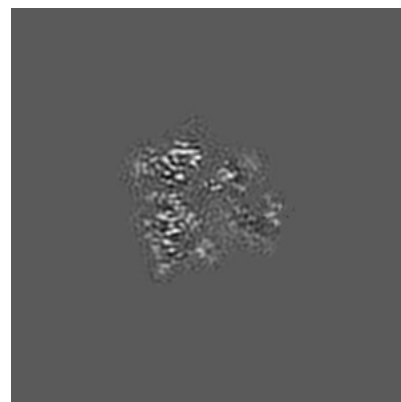
6.2.1 Primary map



X Index: 120



Y Index: 120

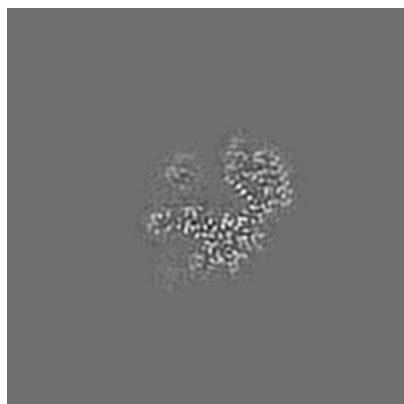


Z Index: 120

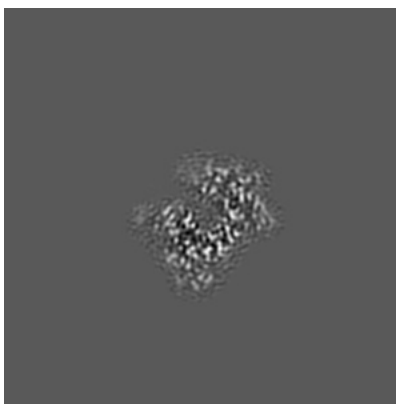
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 109



Y Index: 140

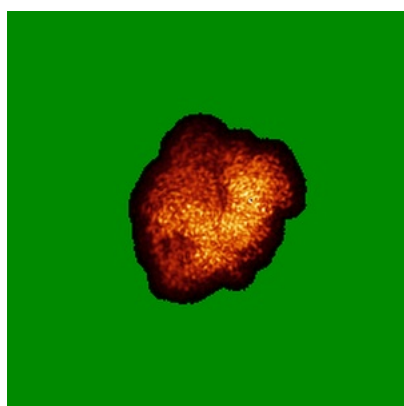


Z Index: 113

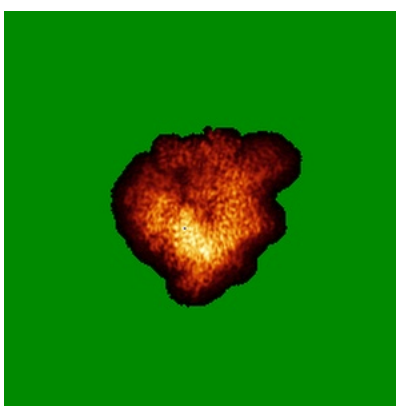
The images above show the largest variance slices of the map in three orthogonal directions.

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

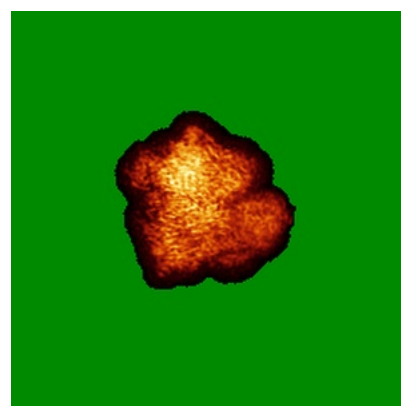
6.4.1 Primary map



X



Y



Z

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 0.01. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

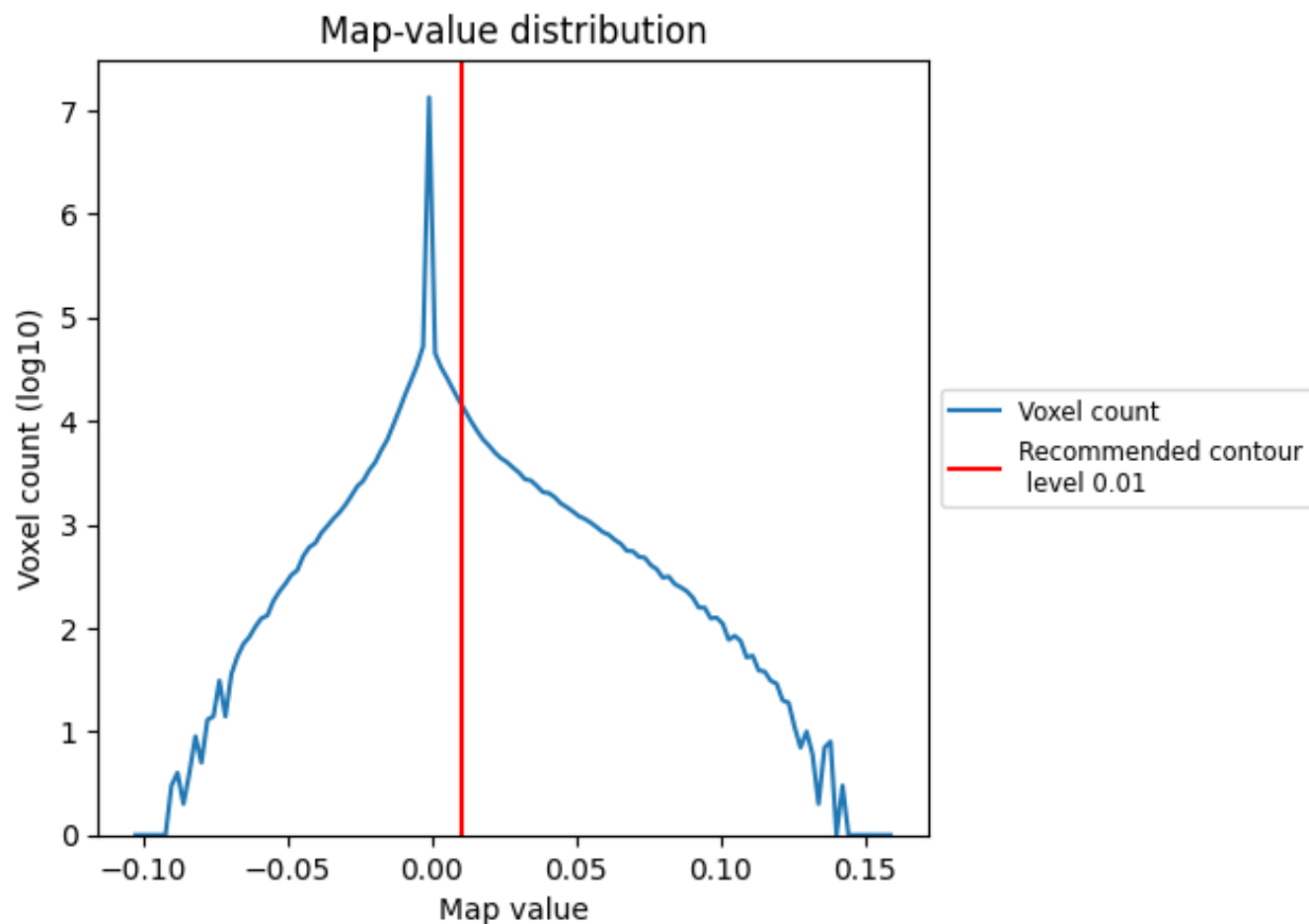
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

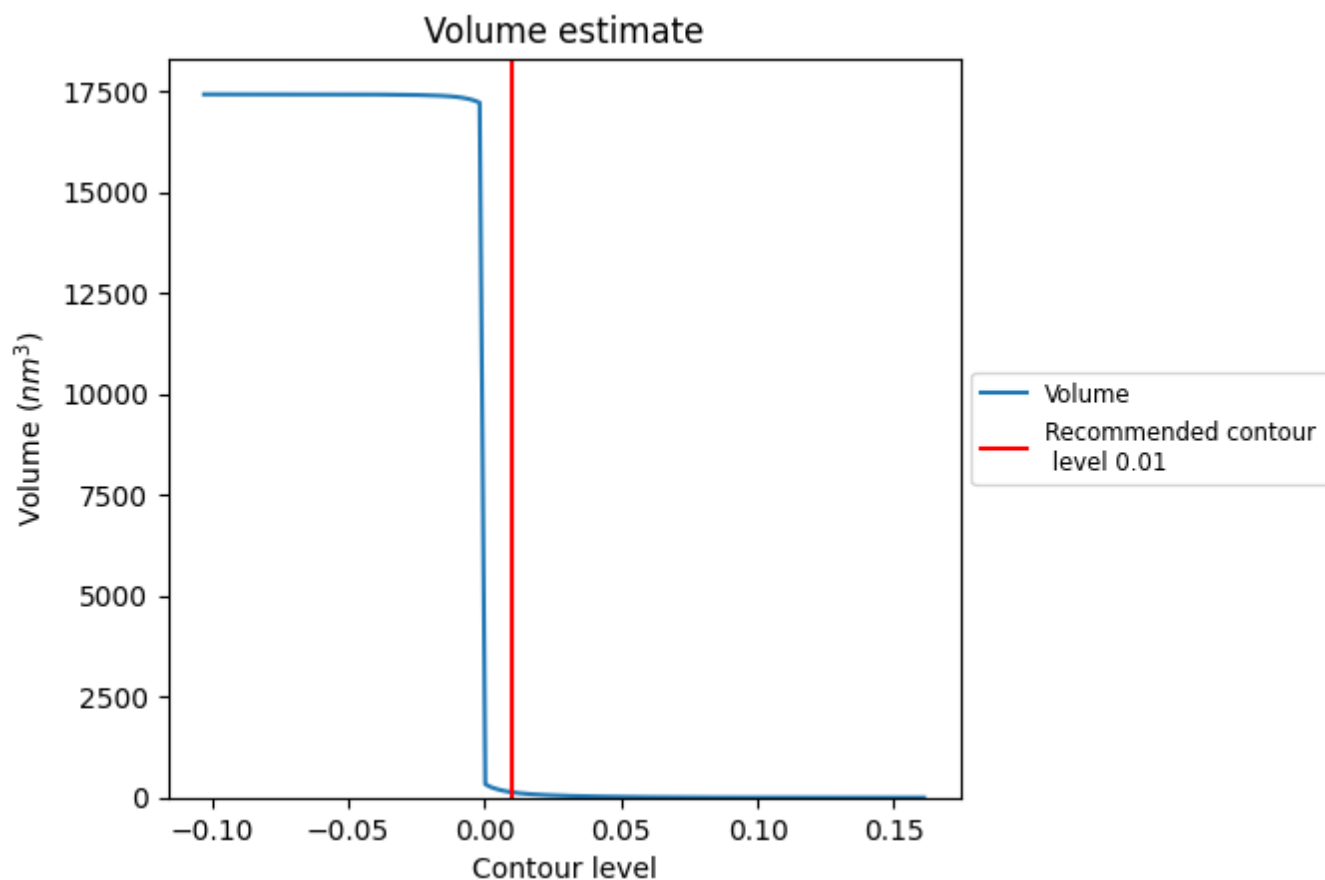
This section contains the results of statistical analysis of the map.

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.

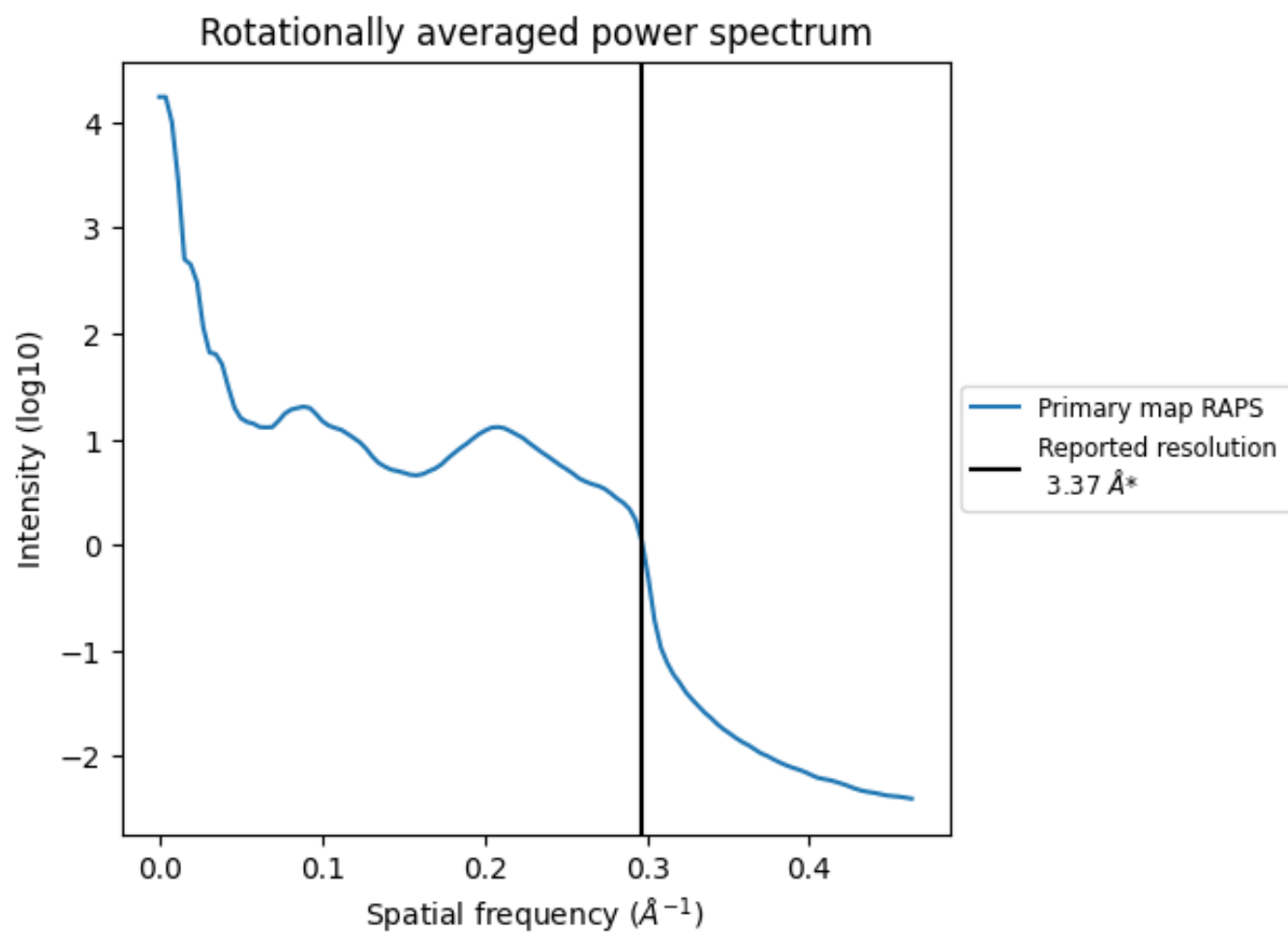
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 133 nm³; this corresponds to an approximate mass of 120 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 ⓘ

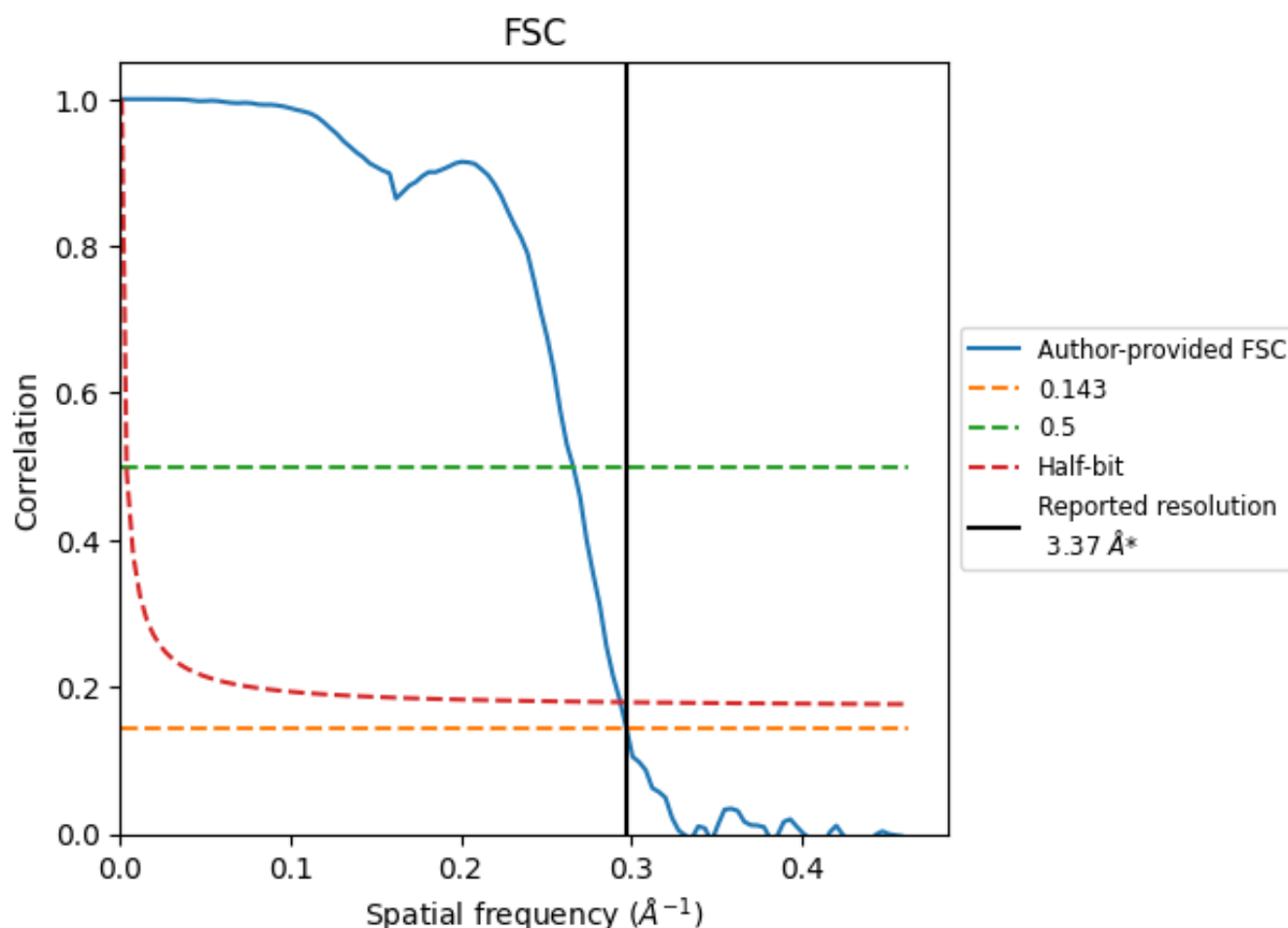


*Reported resolution corresponds to spatial frequency of 0.297 Å⁻¹

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 0.297 \AA^{-1}

8.2 Resolution estimates [i](#)

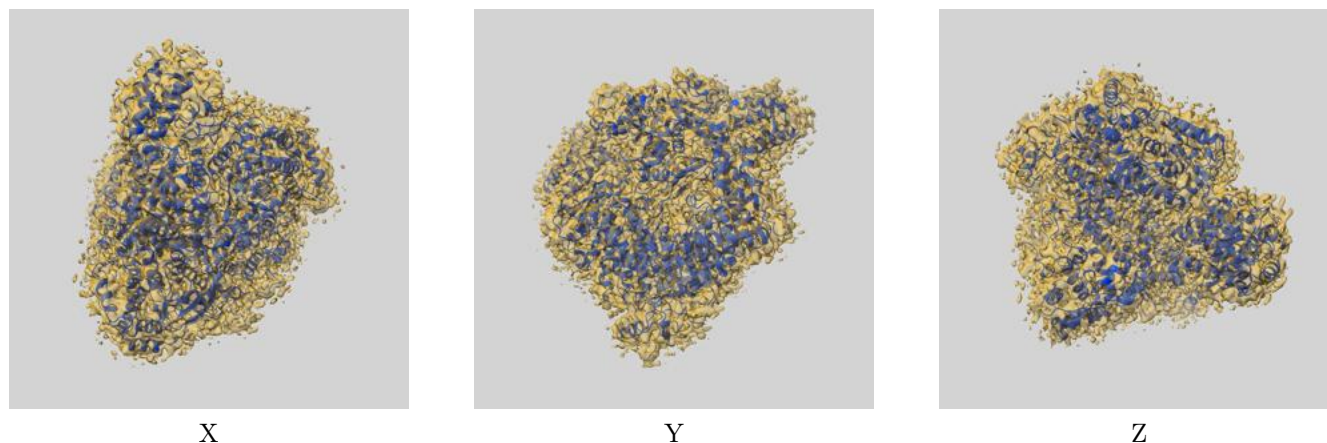
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.37	-	-
Author-provided FSC curve	3.36	3.76	3.40
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

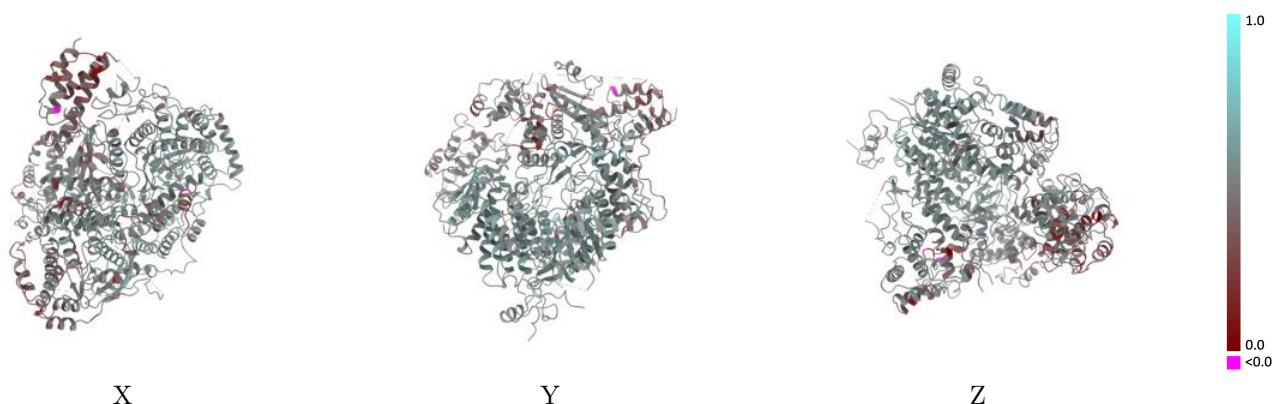
This section contains information regarding the fit between EMDB map EMD-30388 and PDB model 7CKM. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

9.1 Map-model overlay [i](#)



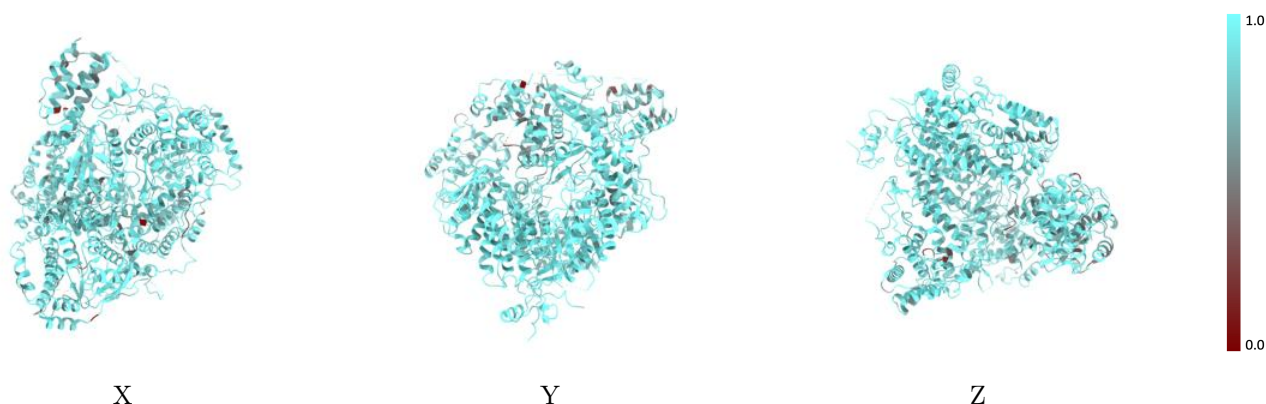
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



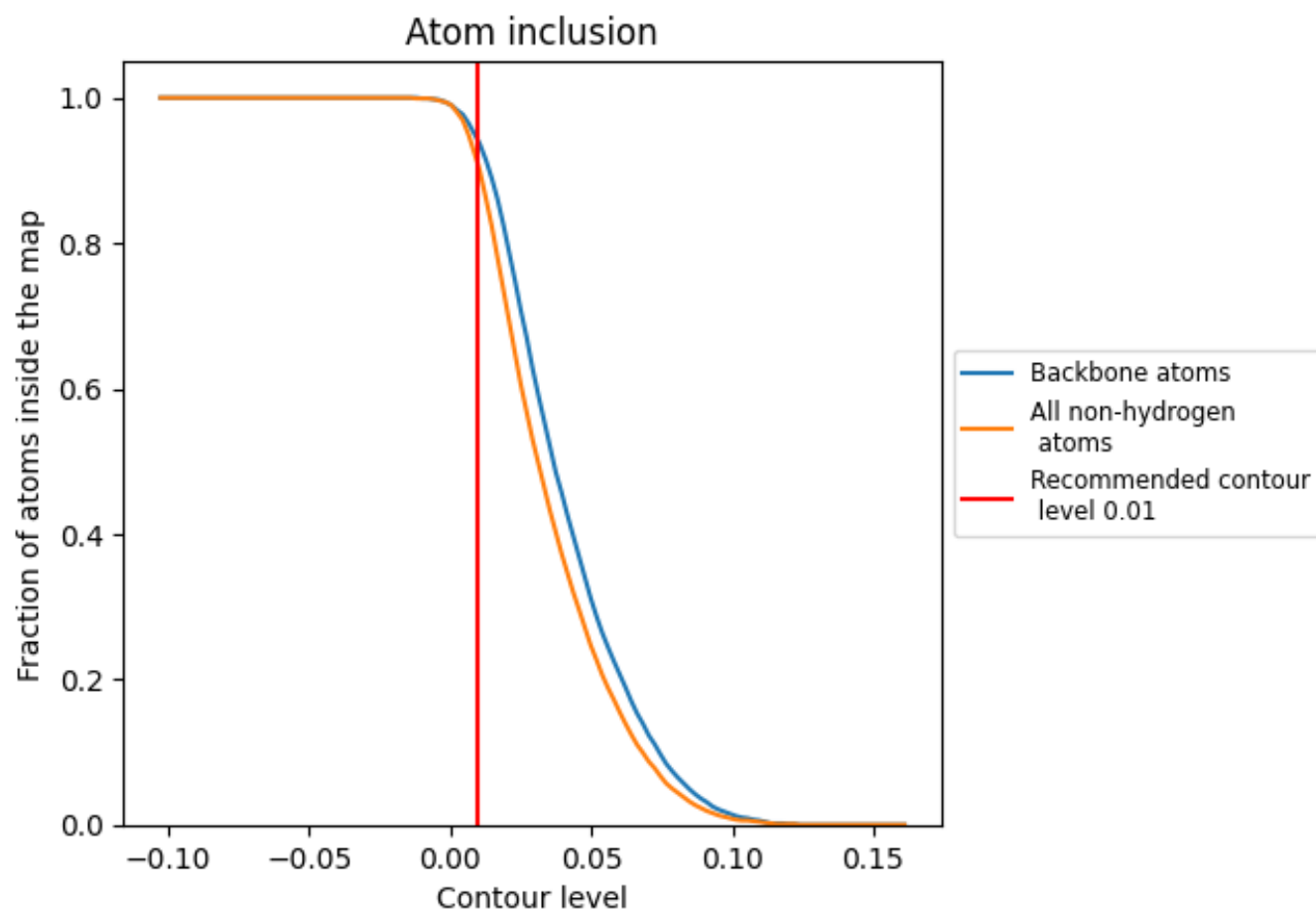
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 91% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.9060	<div></div> 0.4910
A	<div></div> 0.9040	<div></div> 0.4900
B	<div></div> 0.9520	<div></div> 0.5230

