



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 11:56 AM UTC

PDB ID : 7W88 / pdb_00007w88
EMDB ID : EMD-32353
Title : CryoEM structure of open form ZmRDR2 at 3.5 Angstroms resolution
Authors : Du, X.; Yang, Z.; Du, J.
Deposited on : 2021-12-07
Resolution : 3.50 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

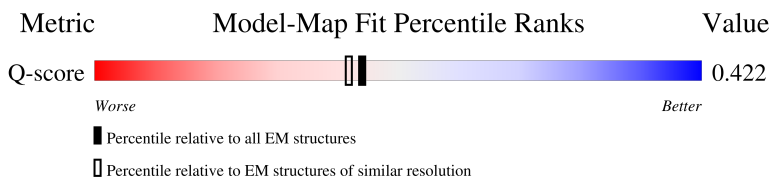
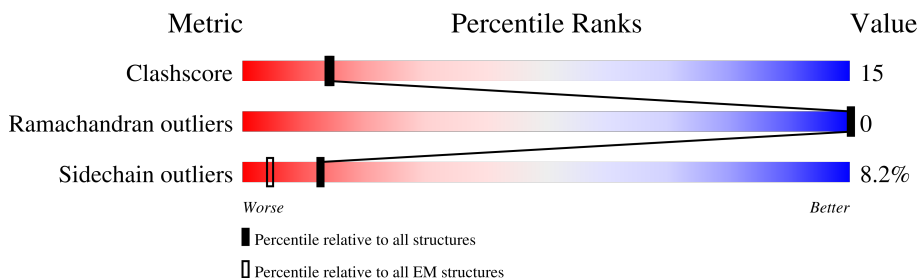
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	13950 (3.00 - 4.00)

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	1131	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7138 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-dependent RNA polymerase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	896	Total	C	N	O	S	0	0
			7138	4541	1241	1311	45		

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	expression tag	UNP Q19VG2
A	-2	MET	-	expression tag	UNP Q19VG2
A	-1	GLY	-	expression tag	UNP Q19VG2
A	0	SER	-	expression tag	UNP Q19VG2
A	962	ALA	GLU	engineered mutation	UNP Q19VG2
A	963	ALA	GLU	engineered mutation	UNP Q19VG2
A	966	ALA	GLU	engineered mutation	UNP Q19VG2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	177075	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.097	Depositor
Minimum map value	-0.054	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.01	Depositor
Map size ($\text{\AA}$)	213.12, 213.12, 213.12	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.74, 0.74, 0.74	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.16	0/7293	0.43	0/9850

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 [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	7138	0	7118	219	0
All	All	7138	0	7118	219	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 15.

All (219) 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:626:VAL:HG23	1:A:717:ILE:HG13	1.68	0.74
1:A:581:GLY:HA2	1:A:610:SER:HB3	1.71	0.72
1:A:521:THR:HG21	1:A:561:VAL:HG22	1.72	0.70
1:A:855:ARG:HH12	1:A:921:PRO:HA	1.57	0.69
1:A:1100:ASP:HA	1:A:1103:LEU:HD12	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ASP:N	1:A:165:ASP:OD1	2.24	0.69
1:A:1081:TYR:HB3	1:A:1096:TRP:HH2	1.58	0.68
1:A:119:ASN:ND2	1:A:234:CYS:O	2.27	0.68
1:A:723:ARG:NH2	1:A:745:THR:O	2.27	0.68
1:A:603:MET:O	1:A:605:LYS:NZ	2.22	0.68
1:A:519:ARG:NH2	1:A:565:ILE:O	2.26	0.67
1:A:642:GLU:OE2	1:A:837:LYS:NZ	2.25	0.67
1:A:746:LYS:HZ2	1:A:750:GLU:HB3	1.61	0.66
1:A:1014:ALA:HB1	1:A:1020:ILE:HG22	1.78	0.65
1:A:627:ASN:OD1	1:A:714:ARG:NH2	2.30	0.65
1:A:223:LEU:HD21	1:A:446:ARG:HD2	1.78	0.64
1:A:492:ARG:HH22	1:A:863:LEU:HD11	1.61	0.64
1:A:735:LEU:HB2	1:A:739:GLN:HB2	1.80	0.63
1:A:115:LEU:HG	1:A:135:VAL:HG21	1.79	0.63
1:A:740:VAL:HG22	1:A:741:TYR:H	1.62	0.63
1:A:777:LYS:NZ	1:A:828:ASP:OD2	2.28	0.62
1:A:656:ASP:HB3	1:A:704:ARG:HG3	1.81	0.62
1:A:649:GLN:HA	1:A:652:MET:HE2	1.81	0.61
1:A:336:LYS:HE3	1:A:1045:ARG:HH22	1.66	0.61
1:A:781:LEU:HD11	1:A:886:HIS:HB2	1.83	0.59
1:A:550:ILE:HD12	1:A:600:ARG:HH11	1.67	0.59
1:A:624:CYS:O	1:A:718:HIS:ND1	2.26	0.59
1:A:785:ASP:OD1	1:A:786:VAL:N	2.35	0.59
1:A:1014:ALA:HA	1:A:1019:GLU:HB2	1.84	0.59
1:A:524:VAL:HG13	1:A:525:SER:H	1.67	0.58
1:A:950:VAL:HB	1:A:954:LEU:HD13	1.85	0.58
1:A:510:ARG:HG2	1:A:605:LYS:HB3	1.85	0.58
1:A:471:GLN:NE2	1:A:476:SER:OG	2.37	0.58
1:A:122:ALA:HB3	1:A:125:GLU:HB2	1.86	0.58
1:A:144:ARG:NH1	1:A:164:GLU:OE2	2.36	0.58
1:A:513:GLN:HG3	1:A:579:ARG:HH21	1.69	0.57
1:A:226:THR:HG21	1:A:230:SER:H	1.69	0.57
1:A:233:GLU:HG2	1:A:279:PRO:HB3	1.87	0.57
1:A:578:ILE:HD12	1:A:587:ILE:HD13	1.86	0.57
1:A:504:VAL:HG23	1:A:861:VAL:HG12	1.87	0.57
1:A:881:ALA:HA	1:A:929:ARG:HD2	1.87	0.56
1:A:317:LYS:O	1:A:317:LYS:NZ	2.32	0.56
1:A:276:ASN:N	1:A:276:ASN:OD1	2.38	0.56
1:A:525:SER:OG	1:A:526:SER:N	2.38	0.56
1:A:1008:LEU:HD23	1:A:1020:ILE:HB	1.88	0.56
1:A:423:PRO:HB3	1:A:875:ILE:HA	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:765:LYS:H	1:A:765:LYS:HZ3	1.54	0.56
1:A:923:GLU:HG2	1:A:924:MET:H	1.69	0.56
1:A:295:PHE:HE2	1:A:388:GLY:H	1.52	0.56
1:A:552:LYS:HE2	1:A:600:ARG:HH22	1.70	0.56
1:A:223:LEU:HD11	1:A:446:ARG:HB2	1.87	0.55
1:A:584:LYS:H	1:A:603:MET:HG3	1.71	0.55
1:A:931:ARG:HB3	1:A:942:ARG:HH22	1.71	0.55
1:A:731:GLU:OE2	1:A:800:ARG:NH2	2.30	0.55
1:A:659:LEU:HD21	1:A:697:LEU:HD13	1.88	0.55
1:A:733:CYS:HB3	1:A:811:ARG:HD2	1.88	0.55
1:A:1095:PRO:HB2	1:A:1096:TRP:HD1	1.72	0.55
1:A:686:GLN:HB3	1:A:1106:LYS:HB3	1.89	0.55
1:A:190:ILE:HB	1:A:222:ALA:HB3	1.89	0.55
1:A:334:MET:HA	1:A:337:LEU:HD23	1.88	0.55
1:A:298:ASN:O	1:A:302:HIS:ND1	2.39	0.54
1:A:603:MET:HE2	1:A:603:MET:H	1.71	0.54
1:A:629:GLU:HG2	1:A:630:ILE:HD12	1.89	0.54
1:A:627:ASN:O	1:A:631:ILE:HG23	2.07	0.54
1:A:628:ARG:H	1:A:628:ARG:HD2	1.72	0.54
1:A:277:VAL:HB	1:A:384:ILE:HG21	1.90	0.54
1:A:241:LEU:HD11	1:A:250:ILE:HD11	1.89	0.54
1:A:518:SER:O	1:A:520:GLN:NE2	2.41	0.54
1:A:331:PHE:O	1:A:335:SER:N	2.38	0.54
1:A:132:CYS:SG	1:A:133:ASP:N	2.81	0.54
1:A:660:THR:OG1	1:A:661:ASN:ND2	2.41	0.54
1:A:658:MET:SD	1:A:658:MET:N	2.78	0.53
1:A:150:LEU:HD11	1:A:225:PHE:CZ	2.44	0.53
1:A:613:LEU:HD12	1:A:613:LEU:H	1.74	0.53
1:A:411:VAL:HG21	1:A:479:MET:HG3	1.90	0.53
1:A:124:ARG:HG2	1:A:260:LEU:HB3	1.91	0.53
1:A:172:CYS:SG	1:A:173:HIS:N	2.82	0.52
1:A:636:THR:O	1:A:896:LYS:NZ	2.42	0.52
1:A:800:ARG:HG3	1:A:802:LEU:HG	1.90	0.52
1:A:1022:THR:OG1	1:A:1026:ARG:NH1	2.42	0.52
1:A:277:VAL:HG23	1:A:280:LEU:HD11	1.90	0.52
1:A:1008:LEU:HB3	1:A:1009:MET:HE2	1.92	0.52
1:A:628:ARG:O	1:A:631:ILE:HG12	2.09	0.52
1:A:937:MET:O	1:A:938:GLU:HG3	2.10	0.52
1:A:882:ILE:HB	1:A:909:HIS:CE1	2.46	0.51
1:A:751:GLN:O	1:A:754:SER:OG	2.28	0.51
1:A:1072:VAL:HG13	1:A:1073:ALA:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:743:ARG:HD2	1:A:768:VAL:HG22	1.92	0.51
1:A:747:ASN:OD1	1:A:747:ASN:N	2.44	0.51
1:A:896:LYS:HD2	1:A:898:ARG:HH22	1.76	0.51
1:A:842:LYS:HD3	1:A:843:VAL:H	1.75	0.51
1:A:618:TRP:HE1	1:A:620:LYS:NZ	2.09	0.51
1:A:869:HIS:HE1	1:A:915:PHE:CD2	2.29	0.51
1:A:941:GLU:HG3	1:A:942:ARG:HG3	1.92	0.51
1:A:469:ALA:O	1:A:473:ARG:NH2	2.35	0.51
1:A:775:ILE:HG21	1:A:808:PHE:HZ	1.75	0.50
1:A:1025:ILE:HG23	1:A:1039:TYR:HE1	1.75	0.50
1:A:471:GLN:OE1	1:A:478:TRP:NE1	2.41	0.50
1:A:555:THR:HA	1:A:589:ILE:HD11	1.92	0.50
1:A:760:CYS:HB3	1:A:768:VAL:HB	1.94	0.50
1:A:940:TRP:HA	1:A:940:TRP:CE3	2.47	0.50
1:A:732:THR:HB	1:A:734:LYS:HE3	1.94	0.50
1:A:626:VAL:O	1:A:627:ASN:ND2	2.45	0.49
1:A:689:GLU:HG2	1:A:691:SER:H	1.77	0.49
1:A:1053:LEU:O	1:A:1057:VAL:HG23	2.13	0.49
1:A:841:GLU:OE1	1:A:842:LYS:N	2.45	0.49
1:A:719:VAL:HG21	1:A:838:LEU:HD22	1.94	0.49
1:A:190:ILE:O	1:A:191:HIS:ND1	2.46	0.49
1:A:525:SER:HG	1:A:527:TYR:HD1	1.57	0.49
1:A:295:PHE:CE2	1:A:386:CYS:HB3	2.48	0.49
1:A:733:CYS:HB3	1:A:811:ARG:HH11	1.77	0.48
1:A:647:MET:SD	1:A:647:MET:N	2.86	0.48
1:A:892:ARG:HG2	1:A:944:MET:SD	2.53	0.48
1:A:746:LYS:N	1:A:765:LYS:O	2.46	0.48
1:A:932:GLU:HB3	1:A:942:ARG:HH12	1.78	0.48
1:A:676:ASP:HB3	1:A:703:HIS:CE1	2.49	0.48
1:A:765:LYS:H	1:A:765:LYS:NZ	2.12	0.48
1:A:1082:HIS:H	1:A:1082:HIS:CD2	2.31	0.48
1:A:999:GLU:O	1:A:1003:GLU:HG2	2.14	0.47
1:A:1017:GLU:HG3	1:A:1093:SER:HB3	1.96	0.47
1:A:628:ARG:HB2	1:A:951:LEU:HD21	1.97	0.47
1:A:644:PHE:HB3	1:A:717:ILE:HD12	1.95	0.47
1:A:283:CYS:HB2	1:A:294:LEU:HD21	1.96	0.47
1:A:551:GLY:HA3	1:A:587:ILE:HG22	1.97	0.47
1:A:662:LYS:NZ	1:A:704:ARG:HH12	2.12	0.47
1:A:1095:PRO:HB2	1:A:1096:TRP:CD1	2.49	0.46
1:A:518:SER:CB	1:A:614:ASN:HB3	2.46	0.46
1:A:704:ARG:O	1:A:708:LEU:HD12	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:923:GLU:HG2	1:A:924:MET:N	2.31	0.46
1:A:245:ALA:HB3	1:A:249:TYR:HB2	1.98	0.46
1:A:520:GLN:CD	1:A:520:GLN:H	2.24	0.46
1:A:632:SER:O	1:A:636:THR:HG23	2.16	0.46
1:A:997:ARG:HG2	1:A:998:TYR:H	1.80	0.46
1:A:507:CYS:O	1:A:511:MET:HB2	2.16	0.46
1:A:695:TYR:HA	1:A:698:MET:HB2	1.97	0.46
1:A:750:GLU:O	1:A:754:SER:OG	2.33	0.45
1:A:725:LEU:HD13	1:A:804:ASP:HA	1.98	0.45
1:A:1058:ARG:O	1:A:1073:ALA:HB2	2.16	0.45
1:A:376:ARG:HB2	1:A:387:LEU:HB2	1.99	0.45
1:A:709:THR:O	1:A:713:THR:HG23	2.16	0.45
1:A:932:GLU:HB3	1:A:942:ARG:NH1	2.32	0.45
1:A:240:LYS:HE2	1:A:240:LYS:HB2	1.77	0.45
1:A:136:ARG:NH1	1:A:138:GLU:OE2	2.50	0.44
1:A:687:GLY:H	1:A:1106:LYS:HG3	1.82	0.44
1:A:465:LEU:HD11	1:A:491:ILE:HB	1.99	0.44
1:A:1045:ARG:HD2	1:A:1045:ARG:HA	1.90	0.44
1:A:503:SER:OG	1:A:918:THR:O	2.35	0.44
1:A:828:ASP:OD1	1:A:828:ASP:N	2.33	0.44
1:A:1009:MET:HE1	1:A:1014:ALA:HB3	1.99	0.44
1:A:455:PHE:CE1	1:A:462:TYR:HD2	2.36	0.44
1:A:235:SER:O	1:A:235:SER:OG	2.30	0.44
1:A:588:THR:OG1	1:A:728:CYS:SG	2.74	0.44
1:A:316:PHE:O	1:A:320:GLU:HB2	2.17	0.44
1:A:321:ASP:OD1	1:A:321:ASP:N	2.50	0.44
1:A:609:LYS:HD3	1:A:609:LYS:HA	1.77	0.44
1:A:997:ARG:HG2	1:A:998:TYR:N	2.33	0.44
1:A:192:ILE:HG13	1:A:222:ALA:HB2	2.00	0.43
1:A:492:ARG:NH2	1:A:863:LEU:HD11	2.31	0.43
1:A:660:THR:HG23	1:A:661:ASN:H	1.83	0.43
1:A:829:LEU:H	1:A:829:LEU:HD23	1.82	0.43
1:A:1025:ILE:HG23	1:A:1039:TYR:CE1	2.53	0.43
1:A:117:LEU:HB3	1:A:234:CYS:SG	2.58	0.43
1:A:952:GLY:O	1:A:956:ARG:NH1	2.52	0.43
1:A:1082:HIS:O	1:A:1084:ASP:N	2.46	0.43
1:A:523:GLU:HG3	1:A:611:THR:HB	1.99	0.43
1:A:550:ILE:O	1:A:587:ILE:HA	2.19	0.43
1:A:740:VAL:HG22	1:A:741:TYR:N	2.32	0.43
1:A:867:GLN:O	1:A:871:VAL:HG23	2.19	0.43
1:A:518:SER:HB2	1:A:614:ASN:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:926:LEU:HG	1:A:927:ALA:H	1.84	0.43
1:A:1040:PHE:HA	1:A:1043:LYS:HE2	2.00	0.43
1:A:249:TYR:CD2	1:A:250:ILE:HG23	2.54	0.43
1:A:608:SER:OG	1:A:609:LYS:N	2.52	0.43
1:A:243:LYS:HD3	1:A:245:ALA:H	1.84	0.43
1:A:297:LEU:O	1:A:301:VAL:HG23	2.19	0.43
1:A:940:TRP:NE1	1:A:943:PRO:HB3	2.34	0.43
1:A:1082:HIS:CG	1:A:1083:PRO:HD3	2.54	0.42
1:A:725:LEU:HB3	1:A:804:ASP:O	2.18	0.42
1:A:525:SER:HG	1:A:527:TYR:H	1.61	0.42
1:A:439:LEU:H	1:A:439:LEU:HD23	1.84	0.42
1:A:489:GLU:O	1:A:493:ARG:HG2	2.20	0.42
1:A:583:TYR:HB3	1:A:603:MET:HE3	2.01	0.42
1:A:1019:GLU:HG2	1:A:1026:ARG:HG3	2.01	0.42
1:A:524:VAL:HG13	1:A:525:SER:N	2.33	0.42
1:A:578:ILE:HD11	1:A:613:LEU:HG	2.02	0.42
1:A:690:PRO:HB2	1:A:1085:ARG:HD3	2.02	0.42
1:A:740:VAL:HB	1:A:789:LEU:HD12	2.01	0.42
1:A:809:PRO:HG2	1:A:816:HIS:CE1	2.55	0.42
1:A:856:ILE:HG22	1:A:858:ASP:H	1.85	0.42
1:A:174:LEU:HB2	1:A:179:ALA:CB	2.49	0.42
1:A:312:ASN:C	1:A:314:ASP:H	2.28	0.42
1:A:554:SER:OG	1:A:556:ARG:HG2	2.19	0.42
1:A:896:LYS:H	1:A:898:ARG:NH1	2.18	0.42
1:A:174:LEU:HD11	1:A:263:LEU:HA	2.02	0.41
1:A:1040:PHE:HA	1:A:1043:LYS:HG2	2.01	0.41
1:A:120:ARG:HH21	1:A:181:LEU:HD21	1.84	0.41
1:A:551:GLY:HA2	1:A:598:SER:O	2.19	0.41
1:A:702:ALA:HA	1:A:1026:ARG:HH22	1.86	0.41
1:A:175:ASP:OD1	1:A:175:ASP:N	2.52	0.41
1:A:911:MET:SD	1:A:911:MET:N	2.93	0.41
1:A:913:VAL:H	1:A:917:LYS:HE2	1.84	0.41
1:A:1106:LYS:HD2	1:A:1106:LYS:HA	1.81	0.41
1:A:373:ARG:HA	1:A:390:GLU:O	2.21	0.41
1:A:841:GLU:CD	1:A:842:LYS:H	2.28	0.41
1:A:1089:LYS:HE2	1:A:1089:LYS:HB2	1.82	0.41
1:A:725:LEU:H	1:A:725:LEU:HG	1.69	0.41
1:A:1058:ARG:HA	1:A:1058:ARG:HD3	1.77	0.41
1:A:547:SER:O	1:A:602:SER:OG	2.28	0.41
1:A:1018:ASP:O	1:A:1022:THR:OG1	2.32	0.41
1:A:931:ARG:HB3	1:A:942:ARG:NH2	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:VAL:HG21	1:A:478:TRP:CZ3	2.57	0.40
1:A:462:TYR:CD1	1:A:481:ALA:HB2	2.57	0.40
1:A:568:ASP:OD1	1:A:568:ASP:N	2.53	0.40
1:A:294:LEU:O	1:A:298:ASN:ND2	2.46	0.40
1:A:679:THR:O	1:A:683:MET:HB2	2.21	0.40
1:A:1041:GLU:H	1:A:1041:GLU:CD	2.27	0.40
1:A:226:THR:CG2	1:A:230:SER:H	2.32	0.40
1:A:174:LEU:HD13	1:A:174:LEU:HA	1.90	0.40
1:A:399:LYS:HE3	1:A:399:LYS:HB2	1.95	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	880/1131 (78%)	746 (85%)	134 (15%)	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	778/954 (82%)	714 (92%)	64 (8%)	10 34

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	VAL
1	A	140	ILE
1	A	150	LEU
1	A	165	ASP
1	A	174	LEU
1	A	180	ILE
1	A	223	LEU
1	A	237	LEU
1	A	276	ASN
1	A	323	SER
1	A	340	THR
1	A	353	HIS
1	A	384	ILE
1	A	391	GLU
1	A	395	ASN
1	A	397	VAL
1	A	437	LYS
1	A	446	ARG
1	A	472	LEU
1	A	473	ARG
1	A	500	ASP
1	A	528	ASP
1	A	536	GLU
1	A	542	THR
1	A	545	ILE
1	A	553	ILE
1	A	564	LEU
1	A	565	ILE
1	A	567	LEU
1	A	578	ILE
1	A	587	ILE
1	A	603	MET
1	A	606	PHE
1	A	630	ILE
1	A	655	SER
1	A	659	LEU
1	A	660	THR
1	A	666	LEU
1	A	668	VAL
1	A	703	HIS
1	A	706	ASN
1	A	708	LEU

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Mol	Chain	Res	Type
1	A	715	CYS
1	A	717	ILE
1	A	728	CYS
1	A	765	LYS
1	A	803	ILE
1	A	806	VAL
1	A	836	ASP
1	A	841	GLU
1	A	875	ILE
1	A	878	THR
1	A	940	TRP
1	A	942	ARG
1	A	997	ARG
1	A	998	TYR
1	A	1020	ILE
1	A	1021	LEU
1	A	1030	VAL
1	A	1081	TYR
1	A	1084	ASP
1	A	1097	ILE
1	A	1099	CYS
1	A	1101	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
1	A	276	ASN
1	A	310	HIS
1	A	346	GLN
1	A	513	GLN
1	A	661	ASN
1	A	686	GLN
1	A	703	HIS
1	A	778	ASN
1	A	810	GLN
1	A	909	HIS
1	A	1027	ASN
1	A	1082	HIS

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](#)

There are no ligands in this entry.

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.

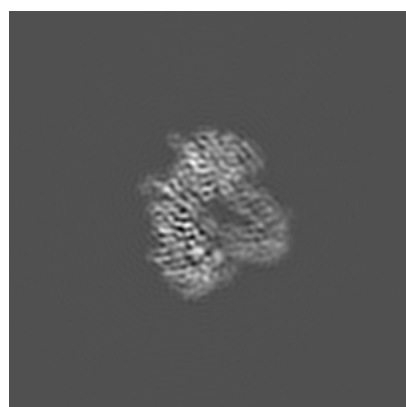
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32353. These allow visual inspection of the internal detail of the map and identification of artifacts.

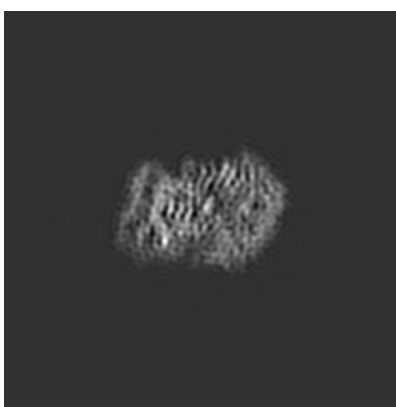
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

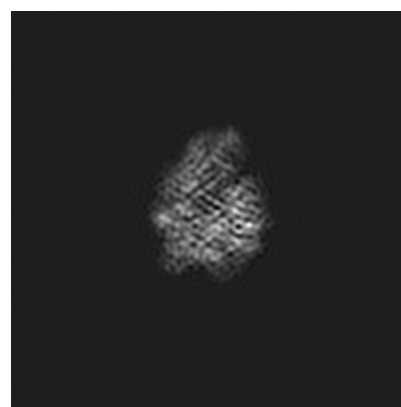
6.1.1 Primary map



X



Y

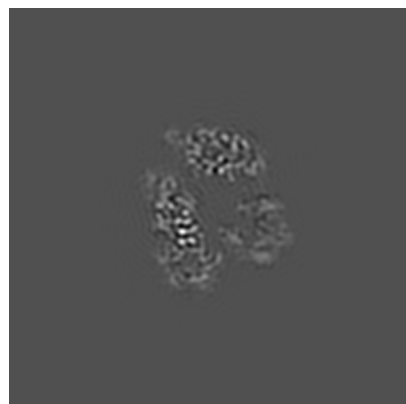


Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 144



Y Index: 144

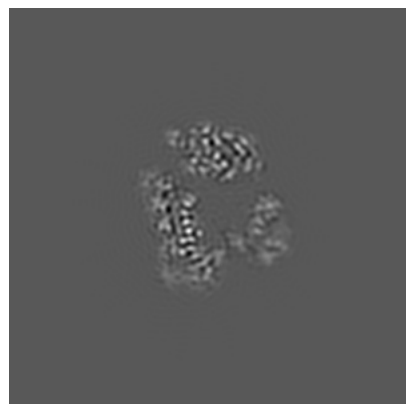


Z Index: 144

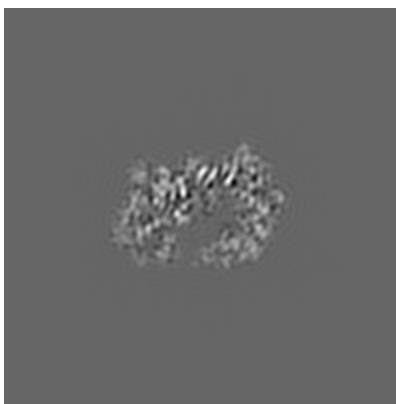
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 148



Y Index: 134

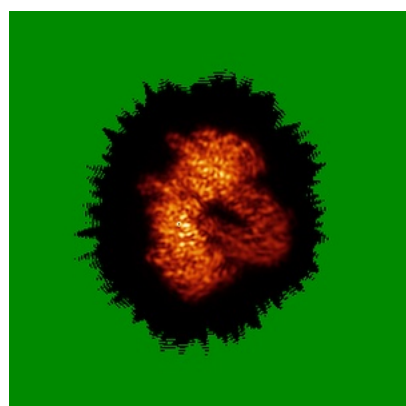


Z Index: 133

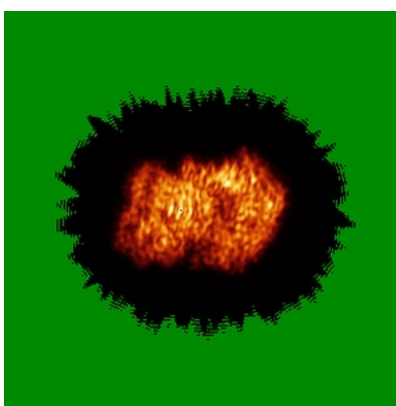
The images above show the largest variance slices of the map in three orthogonal directions.

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

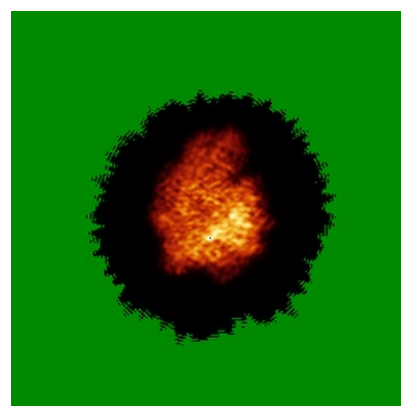
6.4.1 Primary map



X



Y

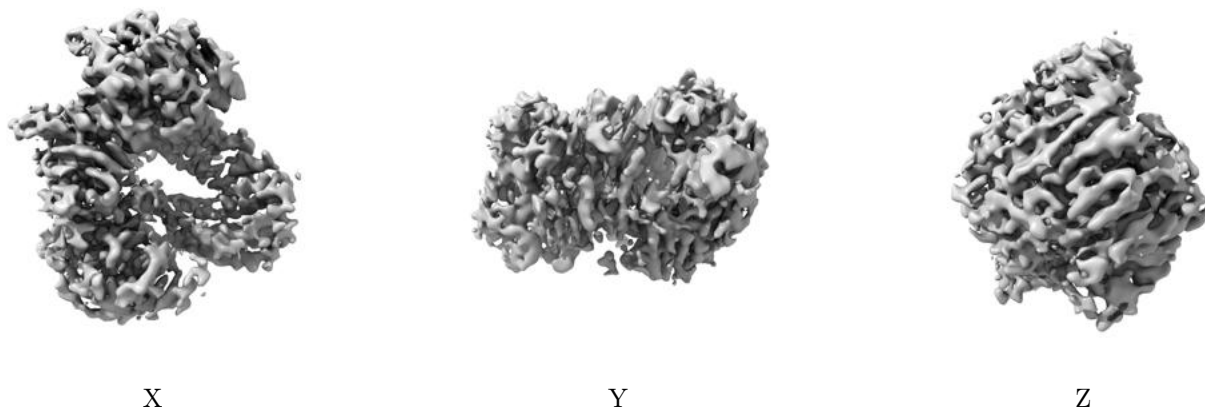


Z

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.

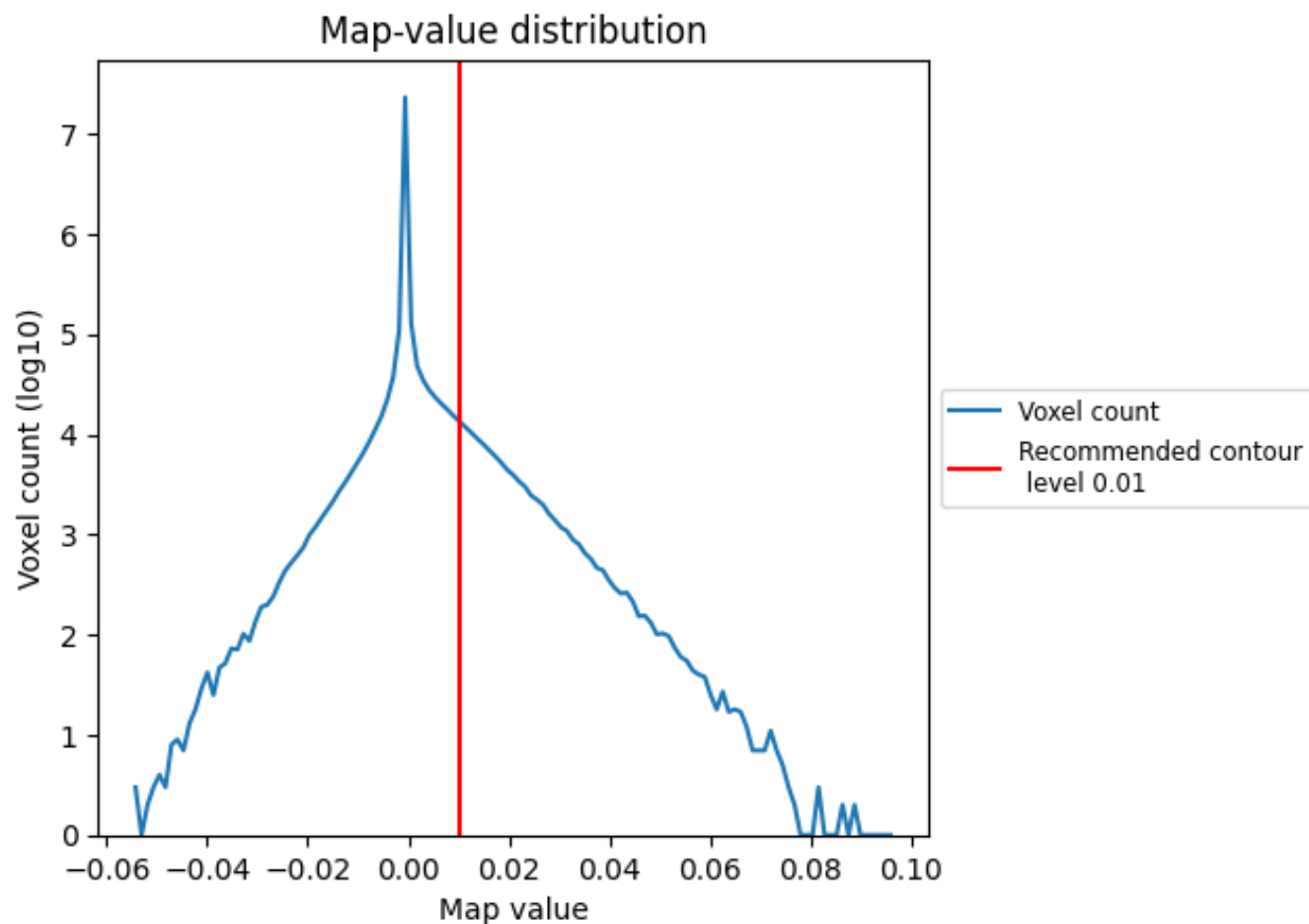
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

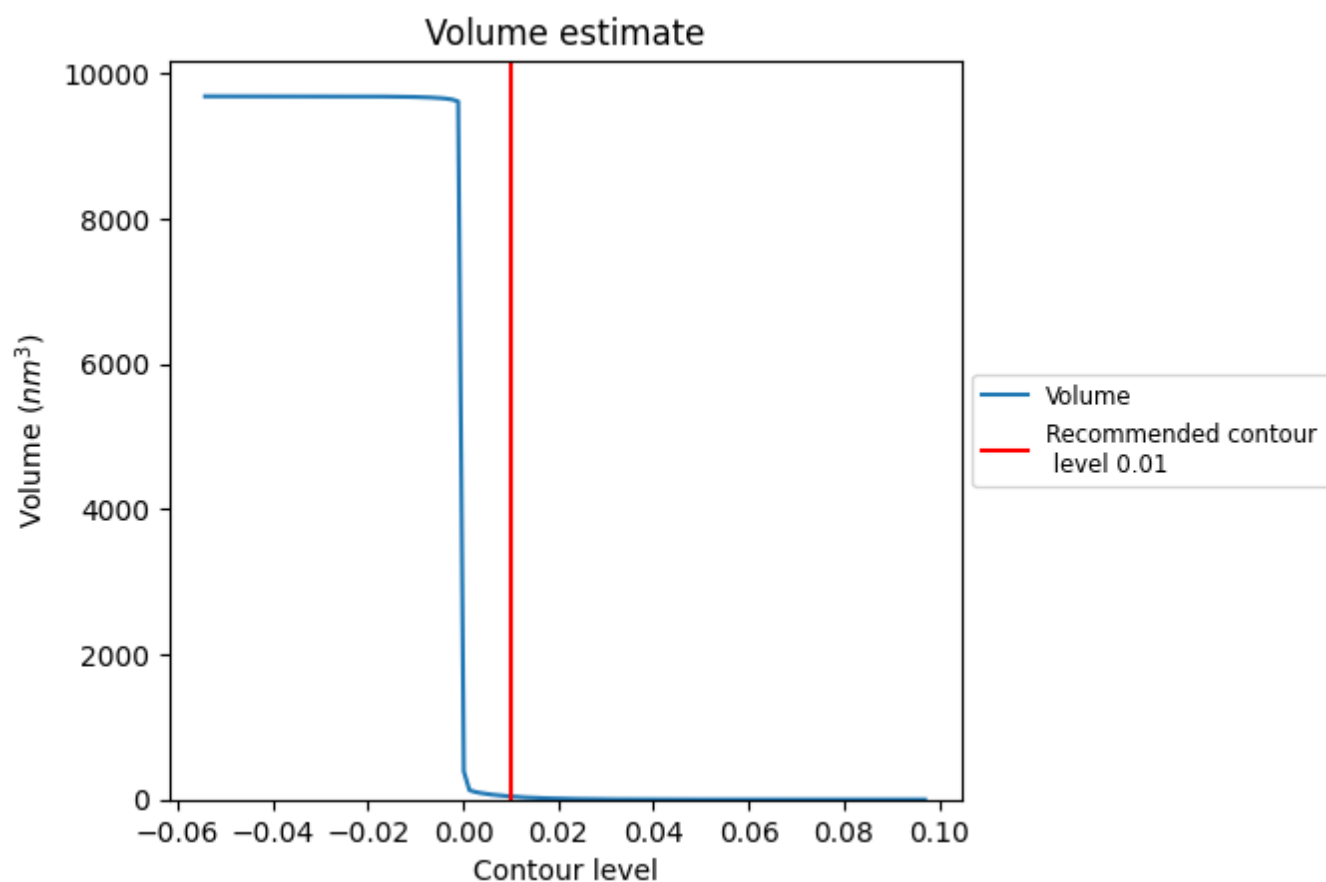
This section contains the results of statistical analysis of the map.

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.

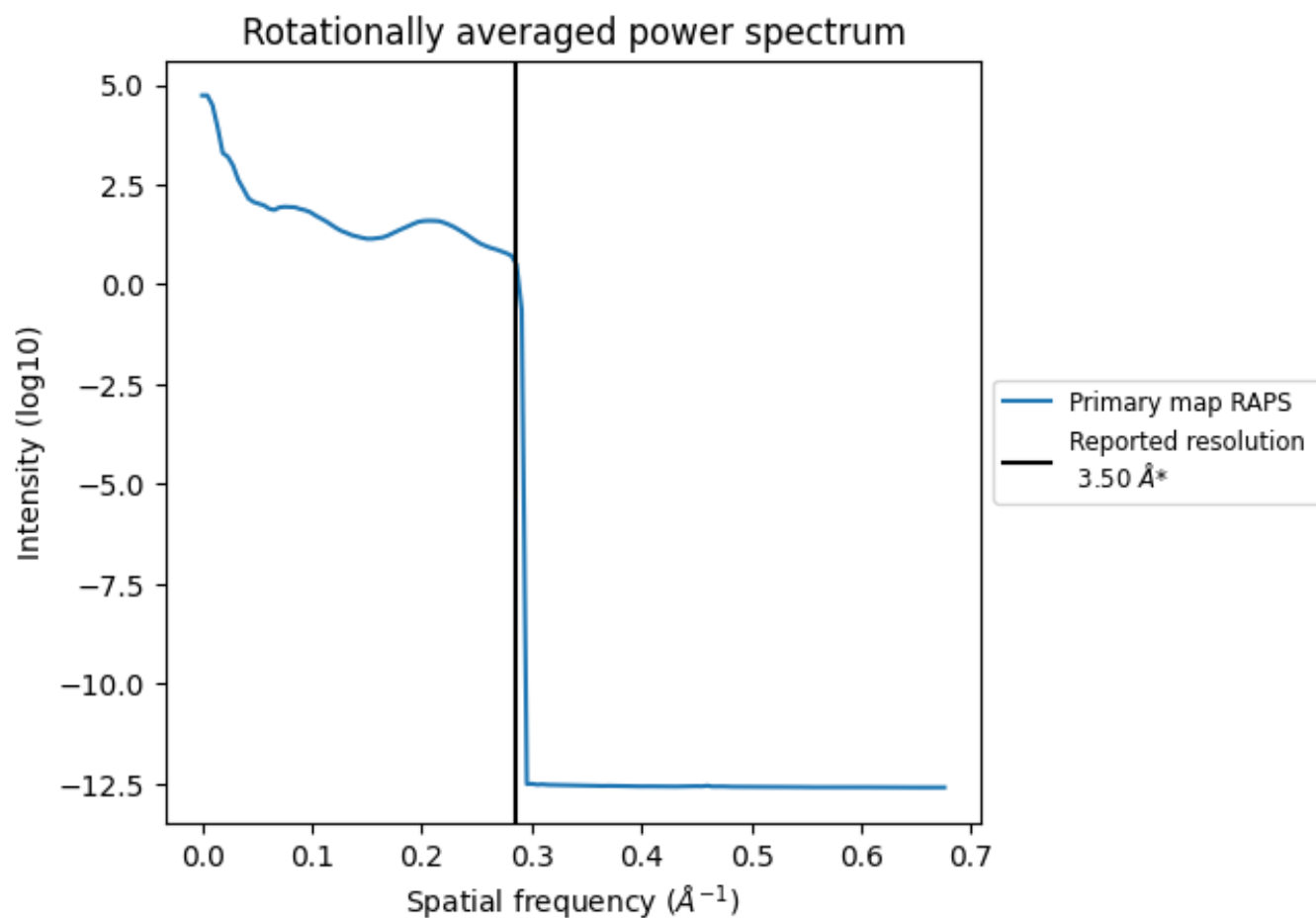
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 43 nm³; this corresponds to an approximate mass of 39 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 0.286 Å⁻¹

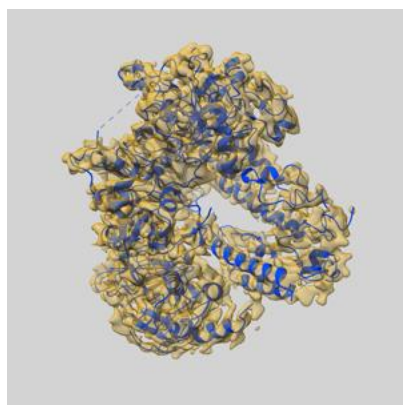
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

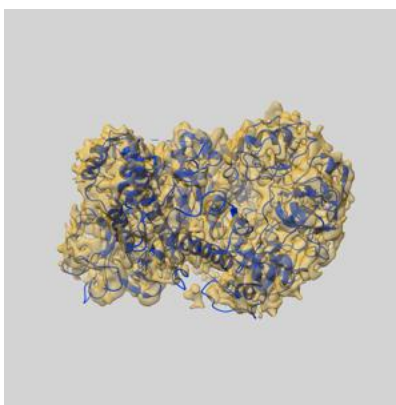
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-32353 and PDB model 7W88. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

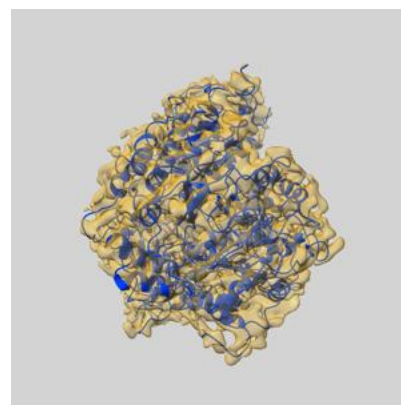
9.1 Map-model overlay [i](#)



X



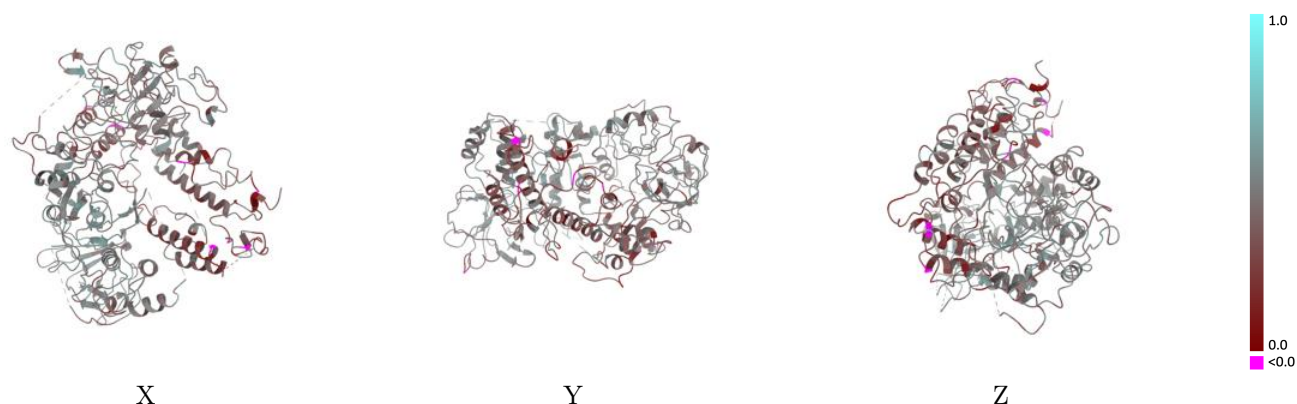
Y



Z

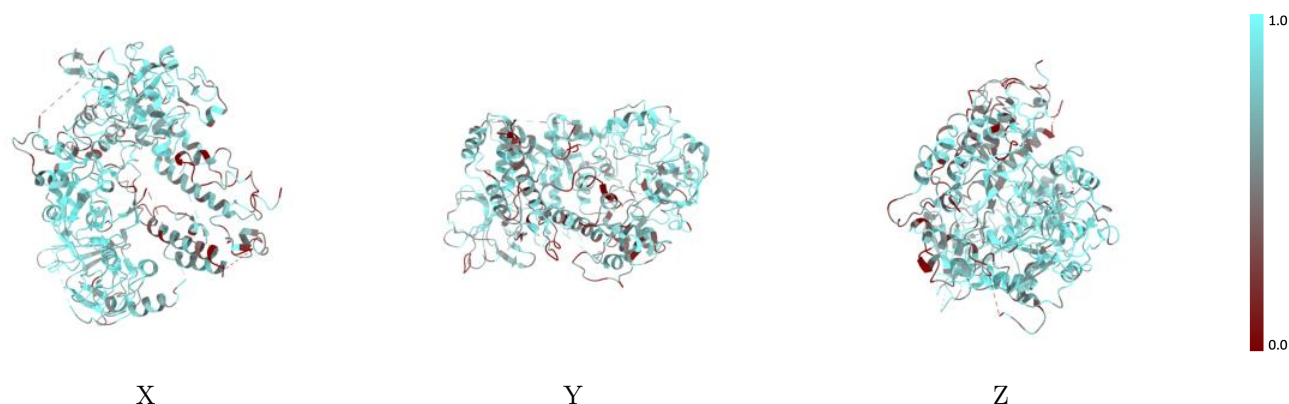
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



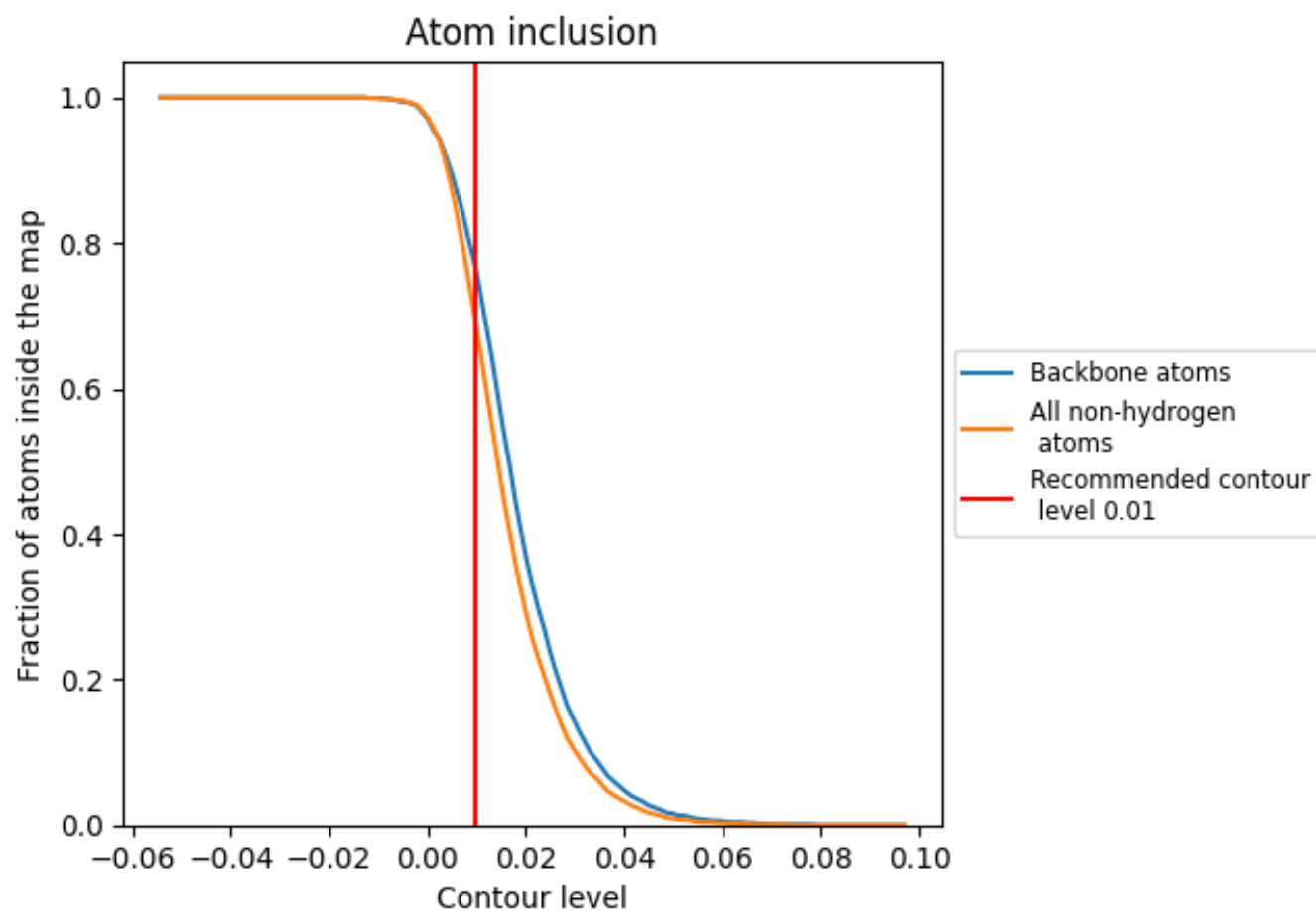
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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, 76% of all backbone atoms, 68% 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	0.6840	0.4220
A	0.6840	0.4220

