



Full wwPDB EM Validation Report ⓘ

Mar 28, 2026 – 09:52 PM UTC

PDB ID : 7VH1 / pdb_00007vh1
EMDB ID : EMD-31983
Title : Cryo-EM structure of Machupo virus dimeric L-Z complex
Authors : Zhang, X.; Ma, J.; Zhang, S.
Deposited on : 2021-09-20
Resolution : 4.20 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

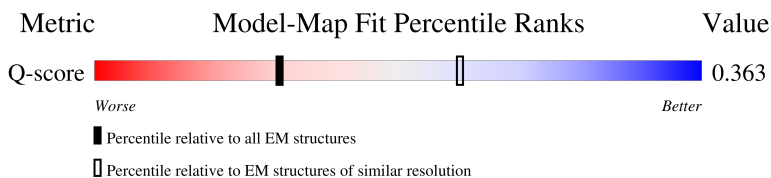
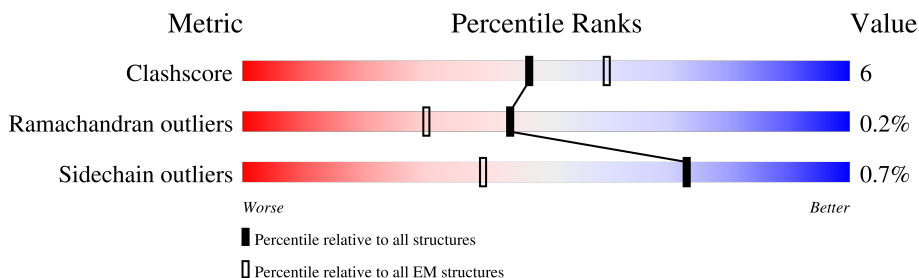
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.20 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	5410 (3.70 - 4.70)

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	2238	
2	B	496	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13156 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	1612	Total	C	N	O	S	0	0
			12749	8144	2129	2389	87		

There are 29 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2210	TRP	-	expression tag	UNP Q6IUF8
A	2211	SER	-	expression tag	UNP Q6IUF8
A	2212	HIS	-	expression tag	UNP Q6IUF8
A	2213	PRO	-	expression tag	UNP Q6IUF8
A	2214	GLN	-	expression tag	UNP Q6IUF8
A	2215	PHE	-	expression tag	UNP Q6IUF8
A	2216	GLU	-	expression tag	UNP Q6IUF8
A	2217	LYS	-	expression tag	UNP Q6IUF8
A	2218	GLY	-	expression tag	UNP Q6IUF8
A	2219	GLY	-	expression tag	UNP Q6IUF8
A	2220	GLY	-	expression tag	UNP Q6IUF8
A	2221	SER	-	expression tag	UNP Q6IUF8
A	2222	GLY	-	expression tag	UNP Q6IUF8
A	2223	GLY	-	expression tag	UNP Q6IUF8
A	2224	GLY	-	expression tag	UNP Q6IUF8
A	2225	SER	-	expression tag	UNP Q6IUF8
A	2226	GLY	-	expression tag	UNP Q6IUF8
A	2227	GLY	-	expression tag	UNP Q6IUF8
A	2228	SER	-	expression tag	UNP Q6IUF8
A	2229	SER	-	expression tag	UNP Q6IUF8
A	2230	ALA	-	expression tag	UNP Q6IUF8
A	2231	TRP	-	expression tag	UNP Q6IUF8
A	2232	SER	-	expression tag	UNP Q6IUF8
A	2233	HIS	-	expression tag	UNP Q6IUF8
A	2234	PRO	-	expression tag	UNP Q6IUF8
A	2235	GLN	-	expression tag	UNP Q6IUF8
A	2236	PHE	-	expression tag	UNP Q6IUF8
A	2237	GLU	-	expression tag	UNP Q6IUF8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	2238	LYS	-	expression tag	UNP Q6IUF8

- Molecule 2 is a protein called Maltose/maltodextrin-binding periplasmic protein,RING finger protein Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	49	Total	C	N	O	S	0	0
			405	254	73	69	9		

There are 37 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-401	MET	-	expression tag	UNP P0AEX9
B	-400	HIS	-	expression tag	UNP P0AEX9
B	-399	HIS	-	expression tag	UNP P0AEX9
B	-398	HIS	-	expression tag	UNP P0AEX9
B	-397	HIS	-	expression tag	UNP P0AEX9
B	-396	HIS	-	expression tag	UNP P0AEX9
B	-395	HIS	-	expression tag	UNP P0AEX9
B	-394	HIS	-	expression tag	UNP P0AEX9
B	-393	HIS	-	expression tag	UNP P0AEX9
B	-26	ASN	-	linker	UNP P0AEX9
B	-25	SER	-	linker	UNP P0AEX9
B	-24	SER	-	linker	UNP P0AEX9
B	-23	SER	-	linker	UNP P0AEX9
B	-22	ASN	-	linker	UNP P0AEX9
B	-21	ASN	-	linker	UNP P0AEX9
B	-20	ASN	-	linker	UNP P0AEX9
B	-19	ASN	-	linker	UNP P0AEX9
B	-18	ASN	-	linker	UNP P0AEX9
B	-17	ASN	-	linker	UNP P0AEX9
B	-16	ASN	-	linker	UNP P0AEX9
B	-15	ASN	-	linker	UNP P0AEX9
B	-14	ASN	-	linker	UNP P0AEX9
B	-13	ASN	-	linker	UNP P0AEX9
B	-12	LEU	-	linker	UNP P0AEX9
B	-11	GLY	-	linker	UNP P0AEX9
B	-10	ILE	-	linker	UNP P0AEX9
B	-9	GLU	-	linker	UNP P0AEX9
B	-8	LEU	-	linker	UNP P0AEX9
B	-7	GLU	-	linker	UNP P0AEX9
B	-6	VAL	-	linker	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	LEU	-	linker	UNP P0AEX9
B	-4	PHE	-	linker	UNP P0AEX9
B	-3	GLN	-	linker	UNP P0AEX9
B	-2	GLY	-	linker	UNP P0AEX9
B	-1	PRO	-	linker	UNP P0AEX9
B	0	GLY	-	linker	UNP P0AEX9
B	1	SER	-	linker	UNP P0AEX9

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	B	2	Total 2	Zn 2	0





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	149686	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TECNAI ARCTICA	Depositor
Voltage (kV)	200	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	0.154	Depositor
Minimum map value	-0.083	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.025	Depositor
Map size (Å)	288.0, 288.0, 288.0	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.6, 1.6, 1.6	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

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	0.44	0/12967	0.76	10/17487 (0.1%)
2	B	0.38	0/417	0.71	0/566
All	All	0.44	0/13384	0.76	10/18053 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	364	ILE	N-CA-C	-7.63	105.33	112.96
1	A	1694	GLY	N-CA-C	6.58	128.77	113.18
1	A	396	ILE	N-CA-C	5.90	121.60	109.34
1	A	699	VAL	N-CA-C	-5.62	105.62	111.58
1	A	1318	TYR	CA-C-N	5.50	130.21	122.46
1	A	1318	TYR	C-N-CA	5.50	130.21	122.46
1	A	395	ASP	CA-C-N	5.47	131.81	121.97
1	A	395	ASP	C-N-CA	5.47	131.81	121.97
1	A	260	LEU	CA-C-N	5.08	131.25	121.54
1	A	260	LEU	C-N-CA	5.08	131.25	121.54

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1693	LEU	Peptide
1	A	1694	GLY	Peptide
1	A	379	GLY	Peptide
1	A	581	PRO	Peptide

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	12749	0	12784	151	0
2	B	405	0	379	8	0
3	B	2	0	0	0	0
All	All	13156	0	13163	158	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 6.

All (158) 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:218:LEU:O	1:A:222:LEU:HB2	1.86	0.74
1:A:382:VAL:H	1:A:385:MET:HE3	1.54	0.71
1:A:285:PRO:HG2	1:A:615:LEU:HD21	1.74	0.70
1:A:971:ASP:HB2	1:A:974:VAL:HG22	1.77	0.67
1:A:1094:LEU:HB2	1:A:1097:TYR:HB2	1.76	0.66
1:A:228:GLY:H	1:A:781:ASN:HD21	1.43	0.66
2:B:41:CYS:SG	2:B:42:CYS:N	2.69	0.66
1:A:1548:ARG:NH1	1:A:1675:SER:O	2.31	0.64
1:A:1552:LEU:HD12	1:A:1630:VAL:HG21	1.80	0.64
1:A:1556:GLY:HA3	1:A:1631:ARG:HD2	1.80	0.64
1:A:774:GLY:H	1:A:1366:ASN:HD22	1.46	0.63
1:A:828:LEU:HG	1:A:1133:VAL:HG21	1.79	0.63
1:A:788:ASN:ND2	1:A:1209:GLU:OE1	2.33	0.61
1:A:1326:THR:HG22	1:A:1328:SER:H	1.65	0.61
2:B:36:ARG:NH2	2:B:38:ASN:OD1	2.34	0.61
1:A:243:LYS:HD3	1:A:1317:LEU:HD23	1.83	0.61
1:A:334:CYS:HB3	1:A:353:LEU:HD12	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:798:LEU:HD21	1:A:802:PRO:HD3	1.81	0.60
1:A:359:ASN:ND2	1:A:437:ARG:O	2.35	0.59
1:A:40:THR:HG21	1:A:1084:SER:HB3	1.85	0.58
1:A:1124:GLN:HE22	1:A:1129:ARG:HH11	1.50	0.58
1:A:1650:CYS:HB2	1:A:1667:SER:HB3	1.86	0.58
1:A:381:ILE:HG22	1:A:386:VAL:HA	1.84	0.57
1:A:456:PRO:HG2	1:A:525:LEU:HB3	1.86	0.57
1:A:388:LEU:HD12	1:A:389:VAL:HG23	1.87	0.57
1:A:1011:LEU:HB3	1:A:1034:GLN:HE22	1.70	0.57
1:A:1495:ARG:NH1	1:A:1682:TYR:OH	2.38	0.57
2:B:72:CYS:SG	2:B:73:HIS:N	2.77	0.57
1:A:573:CYS:SG	1:A:574:ASP:N	2.74	0.56
1:A:586:GLU:HB2	1:A:1438:ARG:HD3	1.87	0.56
1:A:860:LEU:HD21	1:A:1010:LEU:HD21	1.86	0.56
1:A:497:LEU:O	1:A:500:SER:OG	2.22	0.56
1:A:1467:ASP:N	1:A:1467:ASP:OD1	2.39	0.56
1:A:186:ARG:HB2	1:A:1101:SER:HB2	1.87	0.56
1:A:1668:LEU:O	1:A:1672:SER:OG	2.22	0.56
1:A:146:ASN:OD1	1:A:148:ASN:ND2	2.39	0.56
1:A:1192:TRP:HZ3	1:A:1367:LYS:HG2	1.70	0.55
1:A:1641:LYS:O	1:A:1645:SER:N	2.38	0.55
1:A:579:PHE:HB2	1:A:1615:ALA:HB2	1.89	0.54
1:A:1192:TRP:NE1	1:A:1328:SER:O	2.40	0.54
1:A:1402:ALA:O	1:A:1406:ASN:ND2	2.40	0.54
1:A:644:ASN:HB3	1:A:1376:GLY:HA3	1.90	0.54
1:A:678:LEU:O	1:A:692:ARG:NH1	2.41	0.54
2:B:52:CYS:SG	2:B:53:ASN:N	2.65	0.54
1:A:274:GLY:HA3	1:A:672:VAL:HG21	1.90	0.53
1:A:1536:CYS:SG	1:A:1537:SER:N	2.78	0.53
1:A:1464:ASP:OD1	1:A:1464:ASP:N	2.33	0.53
1:A:1161:LEU:HD22	1:A:1297:ASN:HD21	1.74	0.53
1:A:1554:SER:OG	1:A:1555:PHE:N	2.40	0.53
1:A:1739:SER:HB2	1:A:1747:LEU:HD12	1.91	0.53
2:B:38:ASN:ND2	2:B:46:ASP:OD1	2.42	0.52
1:A:385:MET:HG2	1:A:393:VAL:HG13	1.90	0.52
1:A:794:VAL:HG11	1:A:1304:LEU:HD21	1.89	0.52
1:A:984:CYS:HA	1:A:987:VAL:HG12	1.92	0.52
1:A:215:ASP:N	1:A:215:ASP:OD1	2.42	0.51
1:A:201:LEU:HD12	1:A:803:PHE:HA	1.93	0.51
1:A:1780:ASP:HB2	1:A:1788:ILE:HD12	1.94	0.50
1:A:1431:VAL:HG23	1:A:1432:SER:H	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1307:GLU:OE1	1:A:1324:SER:OG	2.28	0.50
1:A:1210:LEU:HD12	1:A:1221:LEU:HD21	1.93	0.50
1:A:368:CYS:SG	1:A:369:GLN:N	2.85	0.49
1:A:678:LEU:HD21	1:A:696:LEU:HD11	1.94	0.49
1:A:1717:TRP:NE1	1:A:1805:GLN:OE1	2.44	0.49
1:A:587:VAL:HG11	1:A:1404:LEU:HD21	1.94	0.49
1:A:479:SER:N	1:A:482:GLU:OE2	2.45	0.49
1:A:106:ARG:NH1	1:A:111:GLU:O	2.45	0.48
1:A:1730:GLN:O	1:A:1734:ASN:ND2	2.46	0.48
1:A:127:LYS:O	1:A:131:GLN:N	2.33	0.48
1:A:1183:LEU:HD13	1:A:1351:TRP:HE1	1.77	0.48
1:A:1502:PHE:HA	1:A:1505:ILE:HG22	1.95	0.48
1:A:1098:SER:OG	1:A:1101:SER:OG	2.27	0.48
1:A:1743:ASN:ND2	1:A:1781:MET:SD	2.86	0.48
1:A:138:VAL:HG11	1:A:845:VAL:HG11	1.95	0.48
1:A:418:SER:HB2	1:A:436:ILE:HG22	1.96	0.48
1:A:1327:SER:O	1:A:1327:SER:OG	2.29	0.48
1:A:712:ASP:HB3	1:A:715:THR:HG22	1.95	0.47
1:A:457:ILE:N	1:A:558:TYR:OH	2.41	0.47
1:A:1716:PHE:HB3	1:A:1804:LYS:HG2	1.97	0.47
1:A:954:VAL:HG13	1:A:957:GLU:HA	1.96	0.47
1:A:1115:LEU:HD13	1:A:1138:THR:HG22	1.97	0.47
1:A:1016:THR:HA	1:A:1019:TYR:HB3	1.97	0.47
1:A:1166:GLU:OE1	1:A:1169:ARG:NH2	2.48	0.47
2:B:67:ARG:NH2	2:B:68:ASN:OD1	2.47	0.46
1:A:343:PHE:HB2	1:A:349:THR:HB	1.97	0.46
1:A:736:VAL:HG12	1:A:737:ASN:H	1.81	0.46
1:A:1202:LEU:HD21	1:A:1229:LYS:HG2	1.97	0.46
1:A:1793:LEU:HD12	1:A:1794:LYS:HG3	1.97	0.46
1:A:274:GLY:O	1:A:276:HIS:ND1	2.48	0.46
1:A:1495:ARG:HD2	1:A:1698:PHE:HA	1.97	0.46
1:A:116:TRP:CD1	1:A:141:VAL:HG21	2.51	0.46
1:A:1087:THR:H	1:A:1090:ALA:HB3	1.80	0.46
1:A:510:LEU:H	1:A:510:LEU:HG	1.58	0.46
1:A:340:LEU:HB3	1:A:342:VAL:HG13	1.97	0.45
1:A:35:PRO:HA	1:A:38:LEU:HB3	1.98	0.45
1:A:50:VAL:HG22	1:A:149:ASN:HD21	1.80	0.45
1:A:584:SER:HB2	1:A:587:VAL:HB	1.98	0.45
1:A:208:LYS:HE3	1:A:210:ASN:HB3	1.96	0.45
1:A:61:GLN:N	1:A:91:PHE:O	2.39	0.45
1:A:204:LEU:HD23	1:A:1223:SER:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1185:CYS:SG	1:A:1333:THR:OG1	2.68	0.45
1:A:81:PRO:HG2	1:A:83:LEU:HG	1.98	0.45
1:A:750:ASP:HA	1:A:753:ILE:HG22	1.98	0.45
1:A:1715:PHE:HB2	2:B:76:TRP:CE2	2.52	0.45
2:B:69:SER:OG	2:B:70:GLU:N	2.50	0.44
1:A:640:MET:HG2	1:A:1379:VAL:HG11	1.99	0.44
1:A:261:THR:HG22	1:A:647:HIS:ND1	2.33	0.44
1:A:124:LEU:HD21	1:A:838:GLU:HG2	2.00	0.44
1:A:646:ILE:HB	1:A:1375:ILE:HG23	2.00	0.44
1:A:1145:VAL:HG11	1:A:1228:LEU:HD21	1.99	0.44
1:A:766:ASP:OD1	1:A:766:ASP:N	2.49	0.43
1:A:104:PHE:HA	1:A:1095:CYS:HB2	2.00	0.43
1:A:1124:GLN:HG2	1:A:1125:VAL:H	1.83	0.43
1:A:969:ASP:HB3	1:A:974:VAL:HG23	2.00	0.43
1:A:1088:ASN:HD22	1:A:1091:LEU:HD21	1.83	0.43
1:A:459:ARG:HA	1:A:522:ASN:HD22	1.83	0.43
1:A:975:LEU:HD12	1:A:976:PRO:HD2	2.00	0.43
1:A:1416:SER:O	1:A:1473:ARG:NH2	2.52	0.43
1:A:130:LEU:HD23	1:A:137:LEU:HD22	1.99	0.43
1:A:775:VAL:HG23	1:A:1365:LEU:HA	2.00	0.43
1:A:455:ARG:HH12	1:A:526:ARG:HG2	1.84	0.43
1:A:718:ILE:HG12	1:A:1242:VAL:HG23	2.01	0.42
1:A:1174:MET:HE2	1:A:1332:ILE:HD13	2.01	0.42
1:A:1183:LEU:HD21	1:A:1375:ILE:HD11	2.00	0.42
1:A:108:ASN:OD1	1:A:108:ASN:N	2.50	0.42
1:A:532:ARG:NH1	1:A:534:TYR:OH	2.51	0.42
1:A:740:LEU:HD21	1:A:1235:VAL:HG11	2.02	0.42
1:A:1515:LEU:HD22	1:A:1530:PHE:HE1	1.85	0.42
1:A:1206:GLN:HA	1:A:1225:LYS:HZ3	1.86	0.41
1:A:1471:GLY:H	1:A:1796:ARG:HH22	1.66	0.41
1:A:1487:ASP:OD1	1:A:1487:ASP:N	2.53	0.41
1:A:355:LEU:HD13	1:A:413:TRP:HZ3	1.84	0.41
1:A:1460:ILE:HG22	1:A:1473:ARG:HH22	1.84	0.41
1:A:1711:ASN:HD22	1:A:1805:GLN:HG2	1.84	0.41
1:A:649:ILE:HD12	1:A:649:ILE:HA	1.94	0.41
1:A:433:ASN:O	1:A:437:ARG:N	2.53	0.41
1:A:501:MET:HE2	1:A:501:MET:HB2	1.92	0.41
1:A:1224:VAL:HA	1:A:1227:ILE:HG22	2.03	0.41
1:A:1718:ALA:HA	1:A:1802:ILE:HA	2.02	0.41
1:A:266:LEU:HD23	1:A:266:LEU:HA	1.89	0.41
1:A:350:LEU:HD21	1:A:392:ARG:NH2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:LEU:HD21	1:A:394:LEU:HB2	2.03	0.41
1:A:455:ARG:H	1:A:524:ARG:HH11	1.67	0.41
1:A:890:LEU:O	1:A:894:TYR:N	2.54	0.41
1:A:1080:LYS:HA	1:A:1083:LYS:HB2	2.03	0.41
1:A:396:ILE:H	1:A:396:ILE:HG13	1.62	0.41
1:A:644:ASN:OD1	1:A:1379:VAL:N	2.46	0.41
1:A:280:ASP:OD1	1:A:280:ASP:N	2.54	0.40
1:A:284:CYS:HA	1:A:285:PRO:HD3	1.89	0.40
1:A:779:LEU:HA	1:A:782:TYR:HD2	1.86	0.40
1:A:848:MET:HA	1:A:851:ARG:HG2	2.04	0.40
1:A:110:ASP:HA	1:A:1136:LEU:HD12	2.03	0.40
1:A:1227:ILE:HG13	1:A:1230:TRP:CE3	2.55	0.40
1:A:1540:GLU:HA	1:A:1543:ARG:HG2	2.02	0.40
1:A:182:GLN:OE1	1:A:186:ARG:NH1	2.54	0.40
1:A:303:VAL:HG22	1:A:485:ARG:HH21	1.87	0.40
1:A:756:LEU:HD23	1:A:756:LEU:HA	1.91	0.40
1:A:588:LEU:HD23	1:A:591:MET:HE2	2.03	0.40
1:A:1516:ILE:H	1:A:1516:ILE:HG13	1.70	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	1582/2238 (71%)	1389 (88%)	189 (12%)	4 (0%)	36	71
2	B	47/496 (10%)	42 (89%)	5 (11%)	0	100	100
All	All	1629/2734 (60%)	1431 (88%)	194 (12%)	4 (0%)	44	77

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	396	ILE
1	A	475	PHE
1	A	383	GLY
1	A	389	VAL

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	1441/2026 (71%)	1430 (99%)	11 (1%)	73	77
2	B	47/412 (11%)	47 (100%)	0	100	100
All	All	1488/2438 (61%)	1477 (99%)	11 (1%)	73	78

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	98	LEU
1	A	394	LEU
1	A	497	LEU
1	A	510	LEU
1	A	704	HIS
1	A	757	LYS
1	A	870	LEU
1	A	1460	ILE
1	A	1560	LEU
1	A	1810	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	34	HIS
1	A	63	ASN
1	A	146	ASN

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Mol	Chain	Res	Type
1	A	148	ASN
1	A	149	ASN
1	A	449	GLN
1	A	472	HIS
1	A	522	ASN
1	A	754	ASN
1	A	771	GLN
1	A	781	ASN
1	A	1034	GLN
1	A	1088	ASN
1	A	1093	ASN
1	A	1111	ASN
1	A	1114	ASN
1	A	1206	GLN
1	A	1297	ASN
1	A	1405	HIS
1	A	1503	ASN
1	A	1689	ASN
2	B	73	HIS

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 2 ligands modelled in this entry, 2 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.

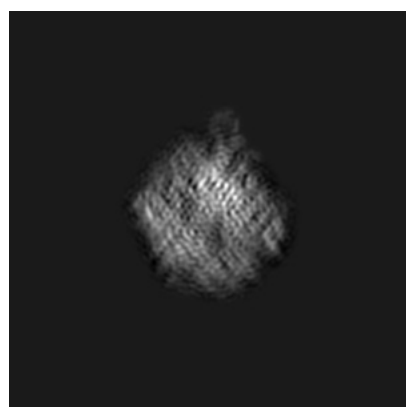
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31983. These allow visual inspection of the internal detail of the map and identification of artifacts.

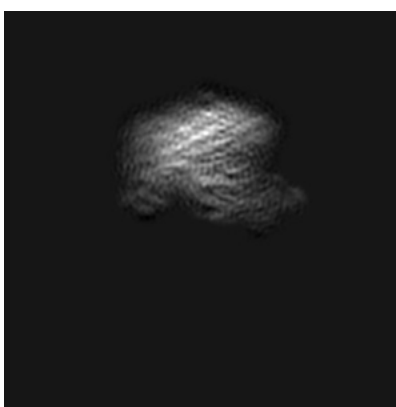
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

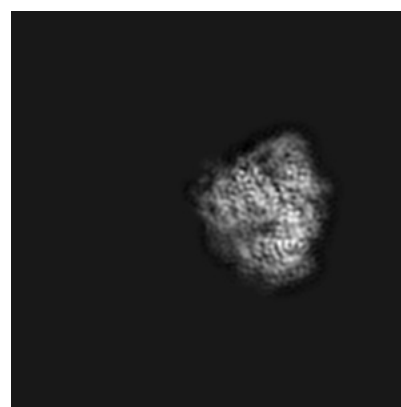
6.1.1 Primary map



X



Y

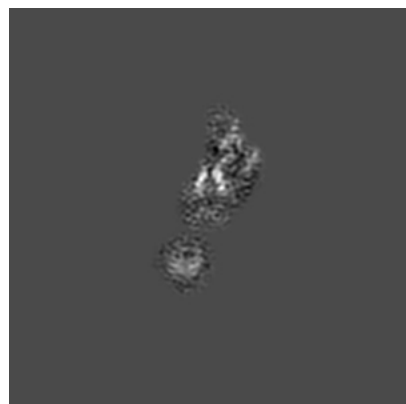


Z

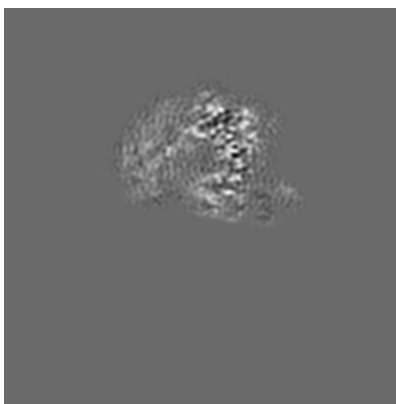
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

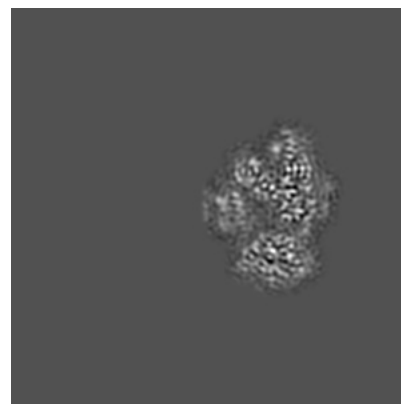
6.2.1 Primary map



X Index: 90



Y Index: 90

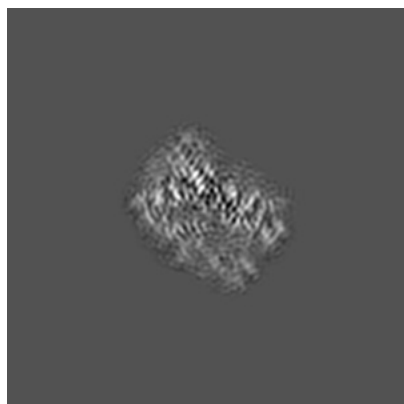


Z Index: 90

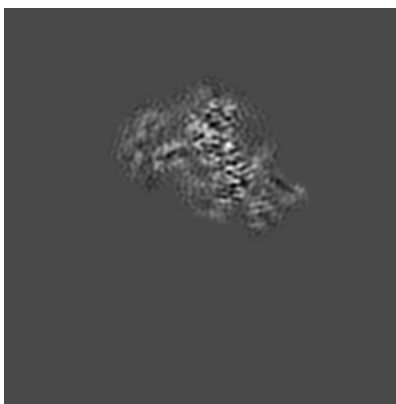
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

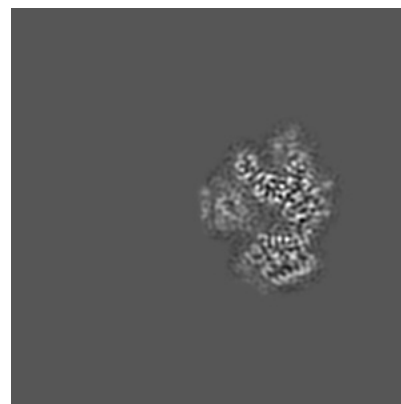
6.3.1 Primary map



X Index: 127



Y Index: 97

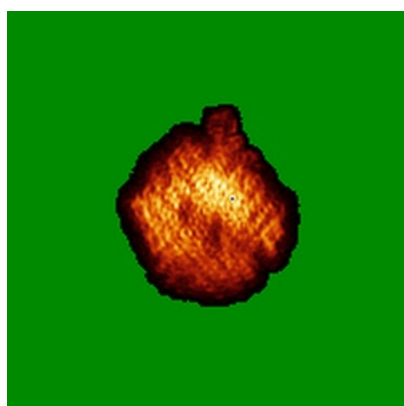


Z Index: 93

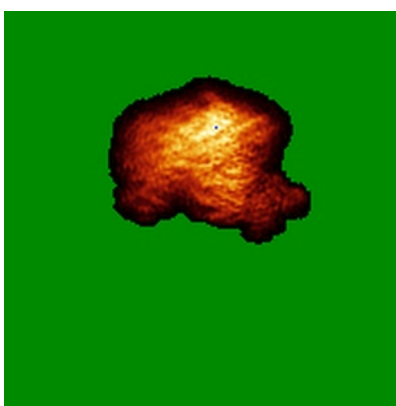
The images above show the largest variance slices of the map in three orthogonal directions.

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

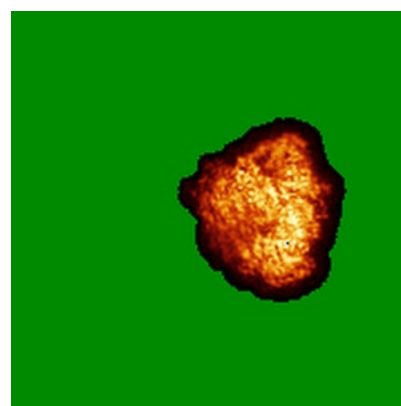
6.4.1 Primary map



X



Y

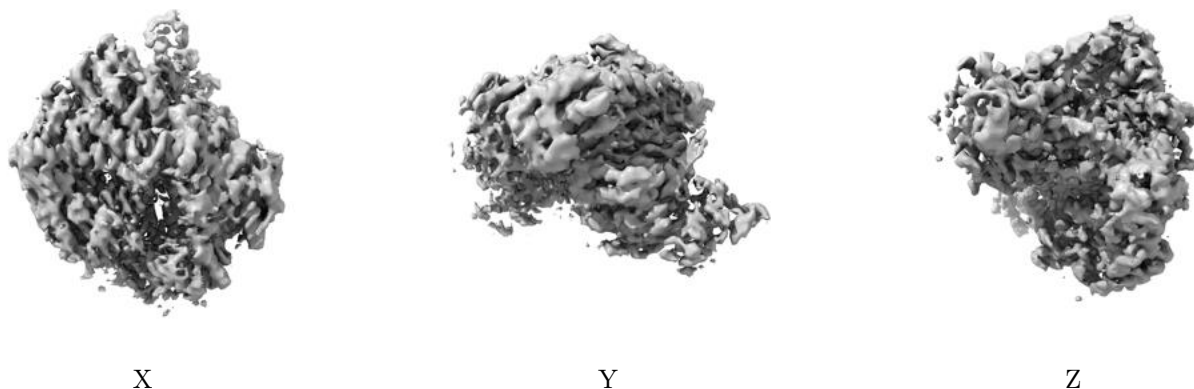


Z

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

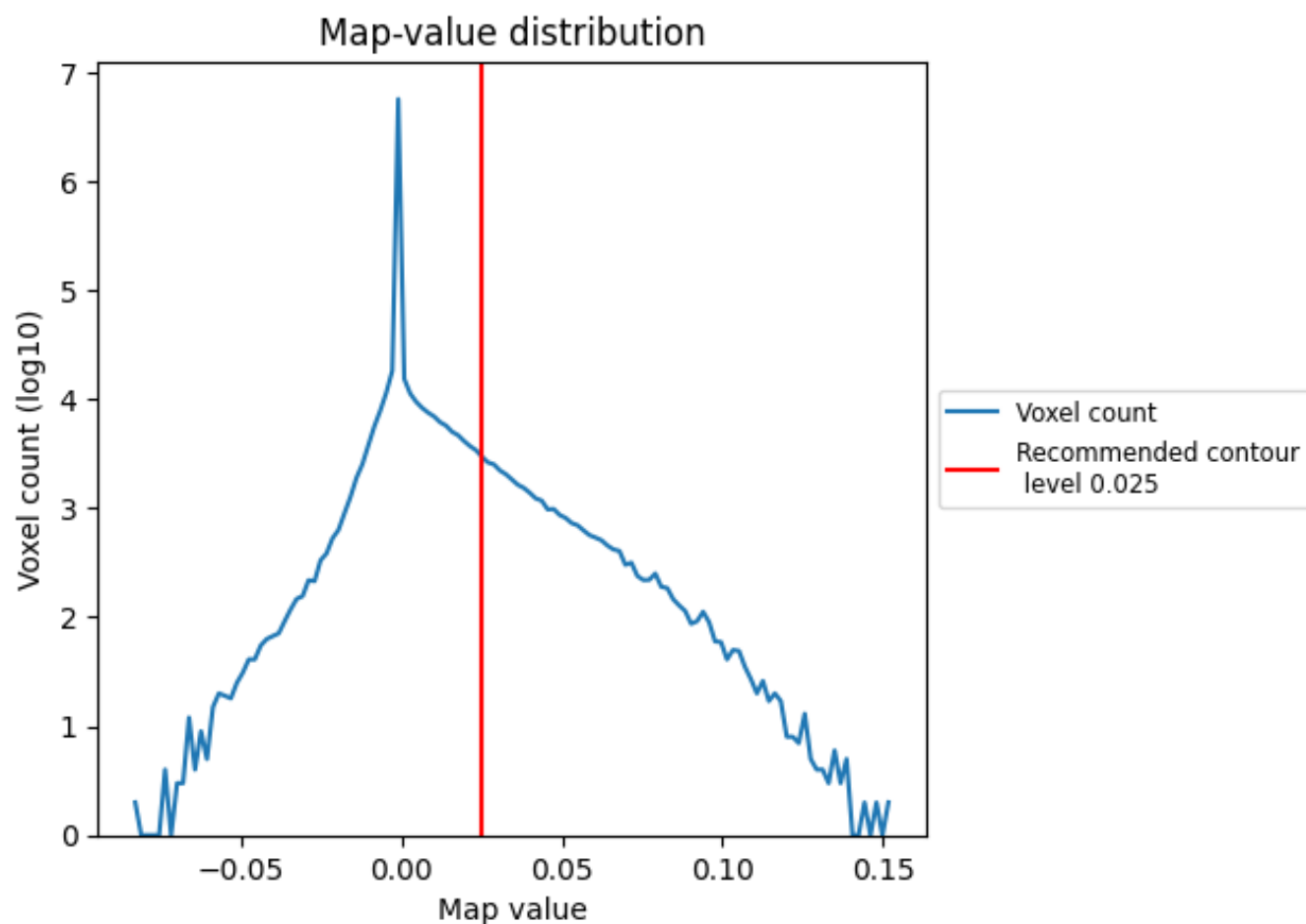
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

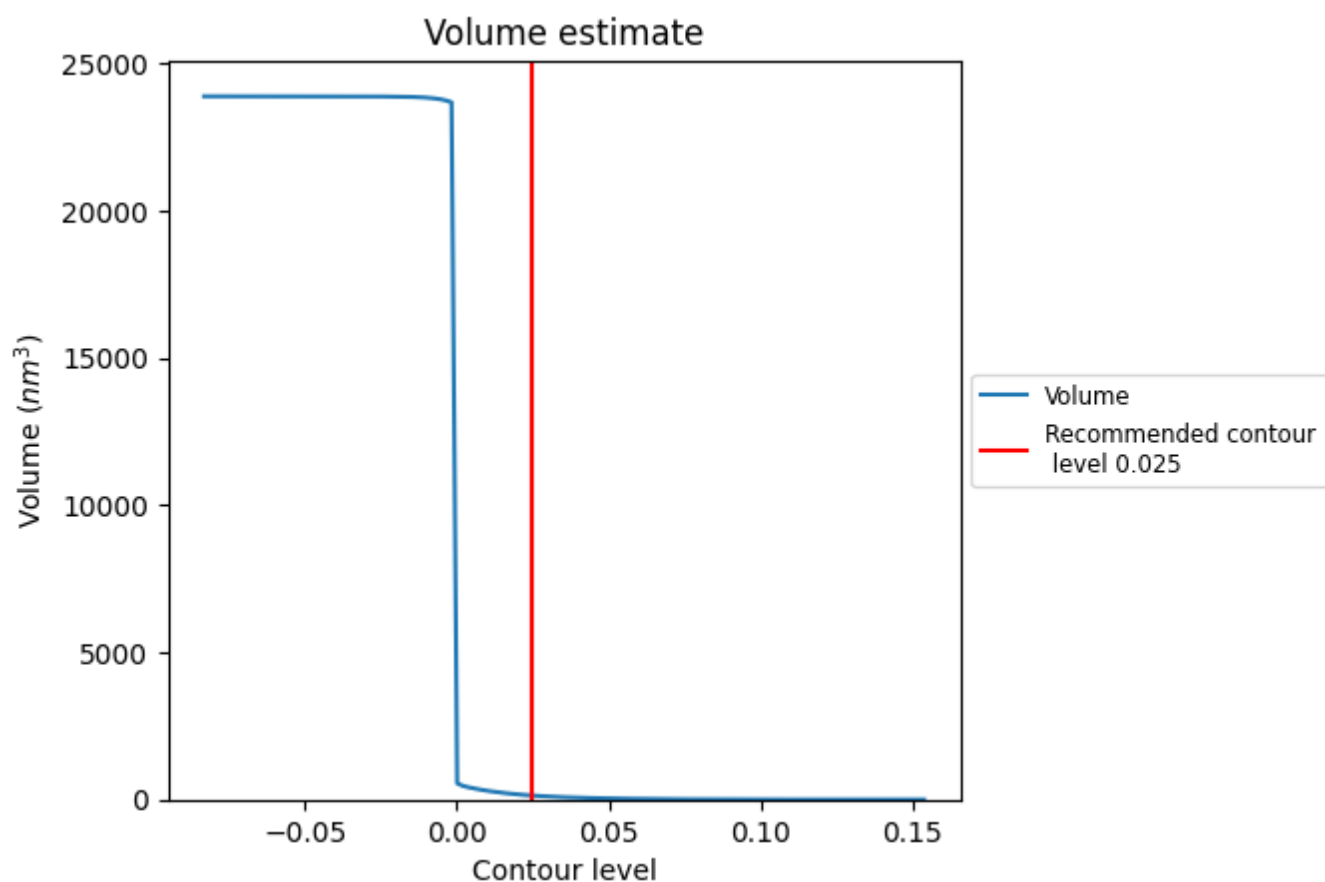
This section contains the results of statistical analysis of the map.

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.

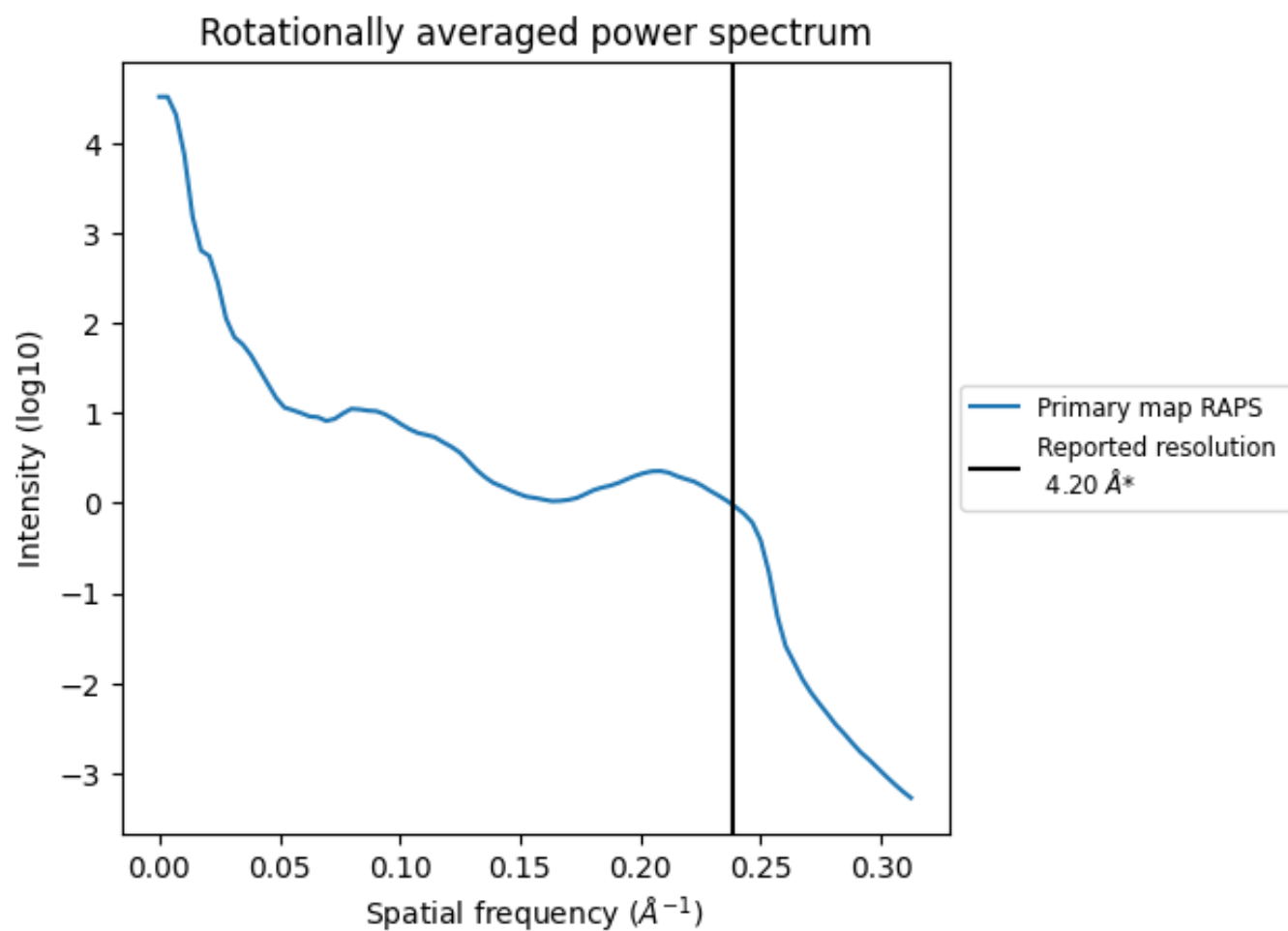
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 135 nm³; this corresponds to an approximate mass of 122 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 0.238 Å⁻¹

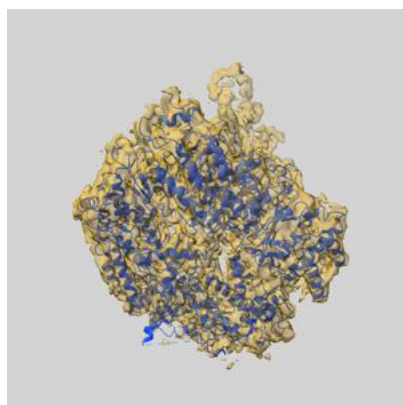
8 Fourier-Shell correlation ⓘ

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

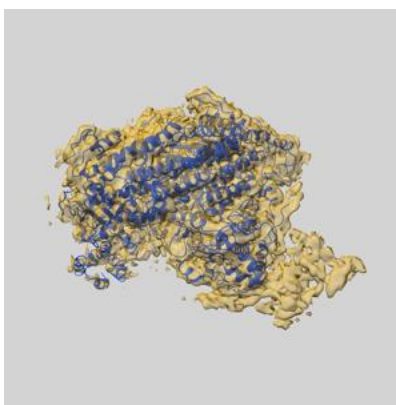
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-31983 and PDB model 7VH1. Per-residue inclusion information can be found in section [3](#) on page [6](#).

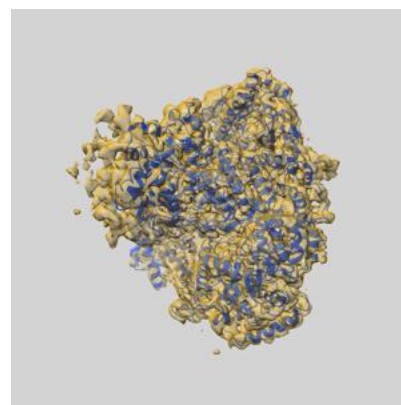
9.1 Map-model overlay [i](#)



X



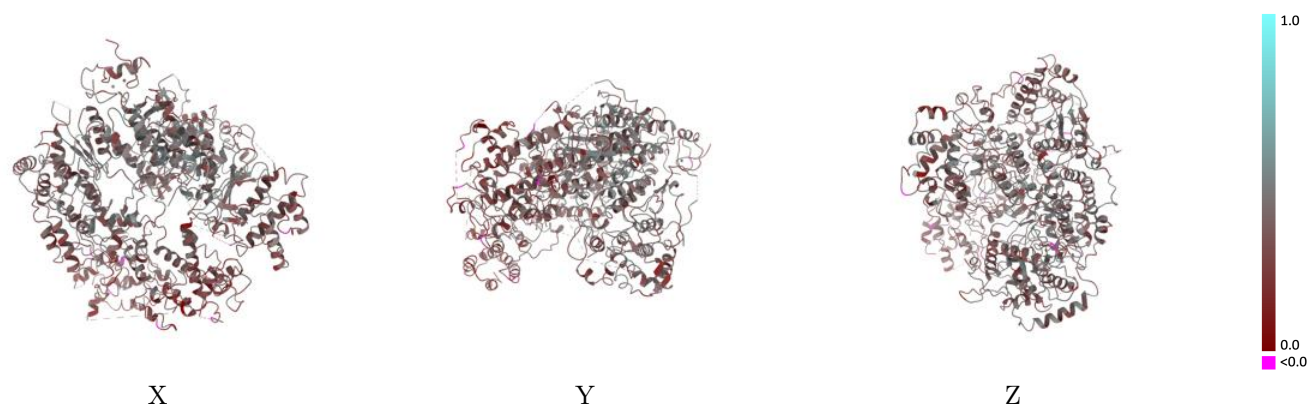
Y



Z

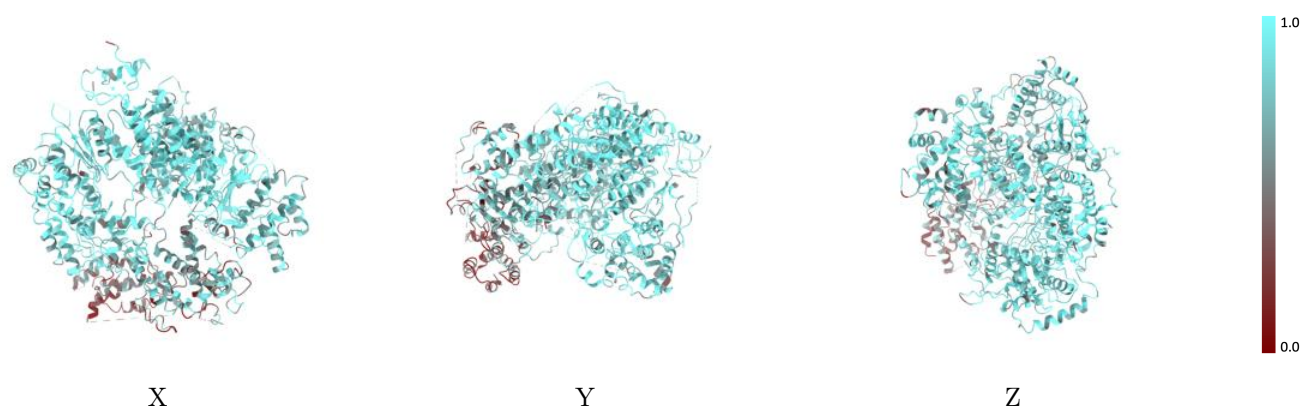
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



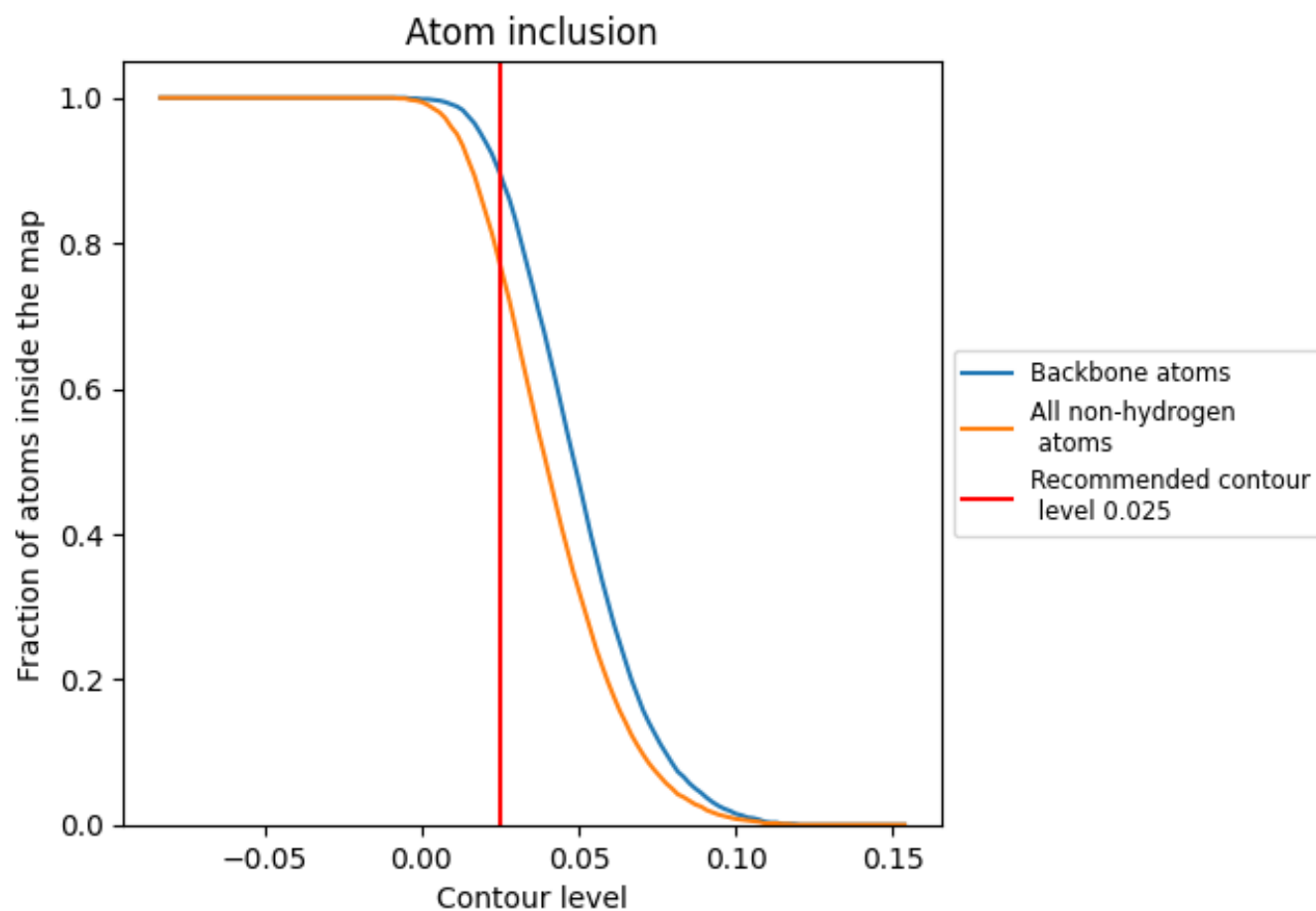
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 77% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7700	0.3630
A	0.7670	0.3630
B	0.8610	0.3470

