



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 03:59 AM UTC

PDB ID : 7VQ7 / pdb_00007vq7
EMDB ID : EMD-32087
Title : The Al-bound AtALMT1 structure at pH 5 (ALMT1Al/pH5)
Authors : Wang, J.Q.
Deposited on : 2021-10-19
Resolution : 3.60 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

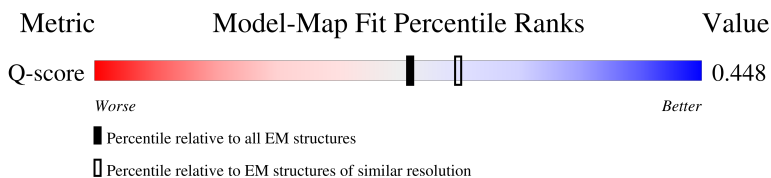
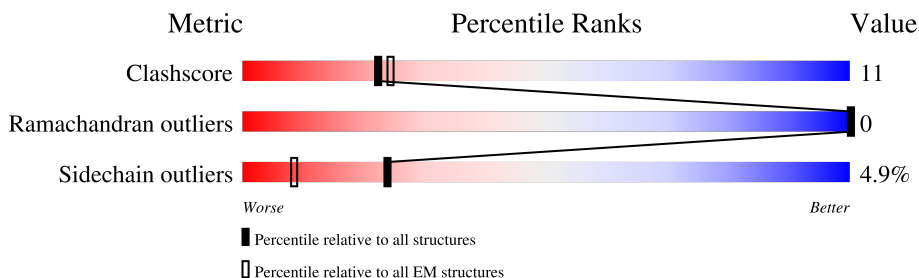
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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	12797 (3.10 - 4.10)

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	509	
1	B	509	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6127 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 Aluminum-activated malate transporter 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	389	Total	C	N	O	S	0	0
			3063	1971	515	560	17		
1	B	389	Total	C	N	O	S	0	0
			3063	1971	515	560	17		

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	494	LEU	-	expression tag	UNP Q9SJE9
A	495	GLU	-	expression tag	UNP Q9SJE9
A	496	GLY	-	expression tag	UNP Q9SJE9
A	497	GLY	-	expression tag	UNP Q9SJE9
A	498	SER	-	expression tag	UNP Q9SJE9
A	499	SER	-	expression tag	UNP Q9SJE9
A	500	GLY	-	expression tag	UNP Q9SJE9
A	501	GLY	-	expression tag	UNP Q9SJE9
A	502	TRP	-	expression tag	UNP Q9SJE9
A	503	SER	-	expression tag	UNP Q9SJE9
A	504	HIS	-	expression tag	UNP Q9SJE9
A	505	PRO	-	expression tag	UNP Q9SJE9
A	506	GLN	-	expression tag	UNP Q9SJE9
A	507	PHE	-	expression tag	UNP Q9SJE9
A	508	GLU	-	expression tag	UNP Q9SJE9
A	509	LYS	-	expression tag	UNP Q9SJE9
B	494	LEU	-	expression tag	UNP Q9SJE9
B	495	GLU	-	expression tag	UNP Q9SJE9
B	496	GLY	-	expression tag	UNP Q9SJE9
B	497	GLY	-	expression tag	UNP Q9SJE9
B	498	SER	-	expression tag	UNP Q9SJE9
B	499	SER	-	expression tag	UNP Q9SJE9
B	500	GLY	-	expression tag	UNP Q9SJE9
B	501	GLY	-	expression tag	UNP Q9SJE9
B	502	TRP	-	expression tag	UNP Q9SJE9
B	503	SER	-	expression tag	UNP Q9SJE9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	504	HIS	-	expression tag	UNP Q9SJE9
B	505	PRO	-	expression tag	UNP Q9SJE9
B	506	GLN	-	expression tag	UNP Q9SJE9
B	507	PHE	-	expression tag	UNP Q9SJE9
B	508	GLU	-	expression tag	UNP Q9SJE9
B	509	LYS	-	expression tag	UNP Q9SJE9

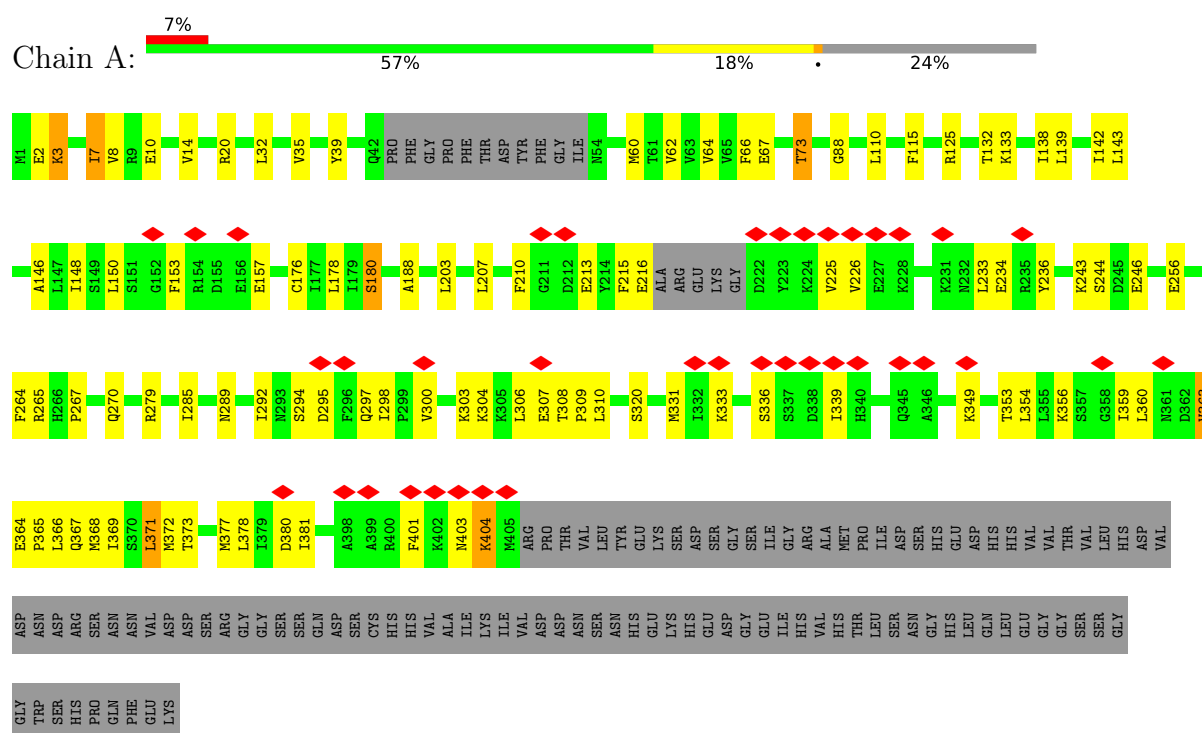
- Molecule 2 is ALUMINUM ION (CCD ID: AL) (formula: Al) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	Al	0
			1	1	

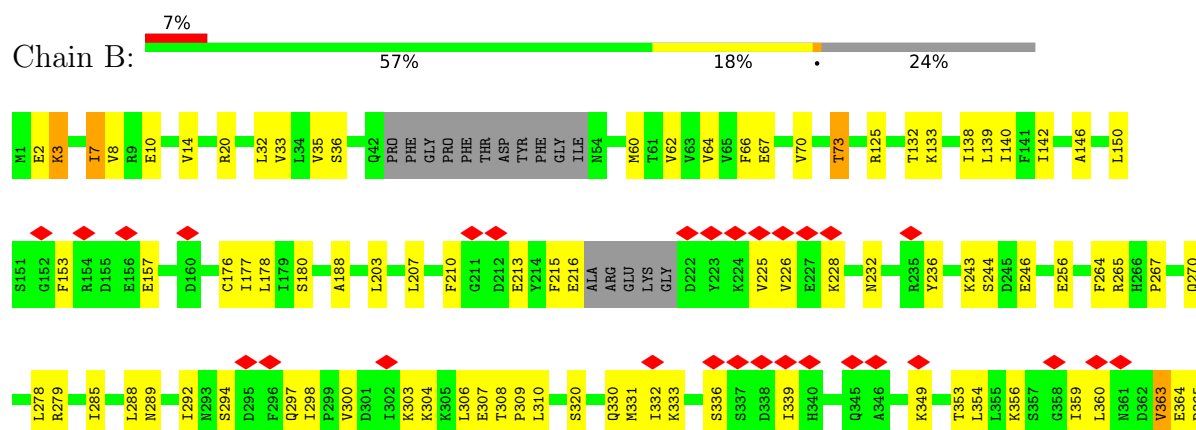
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Aluminum-activated malate transporter 1



• Molecule 1: Aluminum-activated malate transporter 1



SER	ASP	L366
HIS	ARG	Q367
PRO	SER	Q368
GLN	ASN	I369
PHE	ASN	S370
GLU	VAL	L371
LYS	ASP	M372
	ASP	T373
	SER	
	ARG	M377
	GLY	L378
	GLY	I379
	SER	D380
	SER	I381
	GLN	
	ASP	A398
	SER	A399
	CYS	R400
	HIS	F401
	VAL	K402
	ALA	M403
	ILE	K404
	LYS	K404
	ILE	M405
	VAL	ARG
	ASP	PRO
	ASP	THR
	ASN	VAL
	SER	LEU
	ASN	TTR
	HIS	GLU
	GLU	LYS
	LYS	SER
	HIS	ASP
	GLU	SER
	ASP	GLY
	GLY	SER
	GLU	ILE
	ILE	GLY
	HIS	ARG
	VAL	ALA
	HIS	MET
	THR	PRO
	LEU	ILE
	SER	ASP
	ASN	SER
	GLY	HIS
	HIS	GLU
	LEU	ASP
	GLN	HIS
	LEU	VAL
	GLU	VAL
	GLY	VAL
	GLY	THR
	SER	THR
	SER	LEU
	GLY	HIS
	GLY	ASP
	TRP	ASP

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	75357	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 ($e^-/\text{\AA}^2$)	64	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	0.157	Depositor
Minimum map value	-0.107	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.018	Depositor
Map size (Å)	182.52, 182.52, 182.52	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.014, 1.014, 1.014	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AL

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	0.33	0/3120	0.48	0/4211
1	B	0.33	0/3120	0.48	0/4211
All	All	0.33	0/6240	0.48	0/8422

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	3063	0	3126	71	0
1	B	3063	0	3126	74	0
2	A	1	0	0	0	0
All	All	6127	0	6252	141	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 11.

All (141) 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:35:VAL:HG21	1:A:62:VAL:HG21	1.57	0.87
1:B:35:VAL:HG21	1:B:62:VAL:HG21	1.56	0.86
1:B:132:THR:HG23	1:B:133:LYS:HG2	1.70	0.73
1:B:294:SER:O	1:B:297:GLN:NE2	2.22	0.73
1:A:132:THR:HG23	1:A:133:LYS:HG2	1.70	0.73
1:A:294:SER:O	1:A:297:GLN:NE2	2.22	0.71
1:B:210:PHE:HE2	1:B:378:LEU:HD11	1.55	0.71
1:B:353:THR:HA	1:B:356:LYS:HE2	1.73	0.71
1:A:210:PHE:HE2	1:A:378:LEU:HD11	1.56	0.69
1:A:353:THR:HA	1:A:356:LYS:HE2	1.73	0.69
1:A:67:GLU:H	1:A:73:THR:HG23	1.59	0.67
1:B:359:ILE:HG23	1:B:360:LEU:HG	1.78	0.66
1:A:359:ILE:HG23	1:A:360:LEU:HG	1.78	0.65
1:B:67:GLU:H	1:B:73:THR:HG23	1.59	0.64
1:B:2:GLU:OE1	1:B:2:GLU:N	2.30	0.64
1:B:246:GLU:OE2	1:B:279:ARG:NH2	2.30	0.64
1:A:246:GLU:OE2	1:A:279:ARG:NH2	2.30	0.63
1:A:2:GLU:N	1:A:2:GLU:OE1	2.30	0.62
1:A:188:ALA:N	1:A:256:GLU:OE2	2.32	0.62
1:B:2:GLU:HG2	1:B:3:LYS:HD3	1.82	0.62
1:A:2:GLU:HG2	1:A:3:LYS:HD3	1.82	0.61
1:B:188:ALA:N	1:B:256:GLU:OE2	2.32	0.61
1:A:210:PHE:HD1	1:A:292:ILE:HD11	1.66	0.61
1:B:210:PHE:HD1	1:B:292:ILE:HD11	1.66	0.59
1:A:2:GLU:HG2	1:A:3:LYS:H	1.68	0.58
1:A:3:LYS:O	1:A:7:ILE:HG12	2.03	0.58
1:B:3:LYS:O	1:B:7:ILE:HG12	2.03	0.58
1:B:2:GLU:HG2	1:B:3:LYS:H	1.68	0.57
1:A:3:LYS:HD3	1:A:3:LYS:N	2.20	0.56
1:A:176:CYS:O	1:A:180:SER:OG	2.20	0.56
1:A:213:GLU:HA	1:A:216:GLU:HG3	1.86	0.56
1:A:300:VAL:HA	1:A:303:LYS:HE3	1.88	0.56
1:B:213:GLU:HA	1:B:216:GLU:HG3	1.86	0.56
1:B:300:VAL:HA	1:B:303:LYS:HE3	1.88	0.56
1:B:3:LYS:HD3	1:B:3:LYS:N	2.20	0.55
1:B:267:PRO:HB2	1:B:270:GLN:HG3	1.91	0.53
1:B:298:ILE:H	1:B:298:ILE:HD12	1.73	0.52
1:A:298:ILE:H	1:A:298:ILE:HD12	1.73	0.52
1:A:267:PRO:HB2	1:A:270:GLN:HG3	1.91	0.52
1:A:368:MET:O	1:A:372:MET:HG2	2.09	0.52
1:A:349:LYS:O	1:A:353:THR:HG23	2.10	0.51
1:A:146:ALA:O	1:A:150:LEU:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:LYS:HA	1:B:279:ARG:NH2	2.26	0.51
1:B:349:LYS:O	1:B:353:THR:HG23	2.11	0.51
1:B:146:ALA:O	1:B:150:LEU:HB2	2.10	0.51
1:A:210:PHE:CE2	1:A:378:LEU:HD11	2.42	0.51
1:B:368:MET:O	1:B:372:MET:HG2	2.09	0.51
1:A:150:LEU:HD13	1:B:36:SER:HB2	1.93	0.50
1:A:210:PHE:CD1	1:A:292:ILE:HD11	2.46	0.50
1:A:243:LYS:HA	1:A:279:ARG:NH2	2.26	0.50
1:B:364:GLU:HB3	1:B:367:GLN:OE1	2.12	0.50
1:A:236:TYR:OH	1:A:289:ASN:ND2	2.45	0.50
1:A:306:LEU:HB3	1:A:310:LEU:HD12	1.93	0.50
1:A:364:GLU:HB3	1:A:367:GLN:OE1	2.12	0.50
1:B:236:TYR:OH	1:B:289:ASN:ND2	2.45	0.50
1:B:62:VAL:HG13	1:B:176:CYS:SG	2.52	0.50
1:B:306:LEU:HB3	1:B:310:LEU:HD12	1.93	0.49
1:A:62:VAL:HG13	1:A:176:CYS:SG	2.52	0.49
1:B:373:THR:O	1:B:377:MET:HG2	2.13	0.49
1:B:203:LEU:HD22	1:B:285:ILE:HD13	1.94	0.49
1:A:125:ARG:NH1	1:A:139:LEU:HD13	2.28	0.48
1:A:367:GLN:O	1:A:371:LEU:HD12	2.13	0.48
1:B:210:PHE:CD1	1:B:292:ILE:HD11	2.46	0.48
1:B:333:LYS:HB2	1:B:403:ASN:HA	1.95	0.48
1:A:373:THR:O	1:A:377:MET:HG2	2.13	0.48
1:B:3:LYS:HD3	1:B:3:LYS:H	1.78	0.48
1:A:3:LYS:HD3	1:A:3:LYS:H	1.78	0.48
1:A:333:LYS:HB2	1:A:403:ASN:HA	1.95	0.48
1:B:125:ARG:NH1	1:B:139:LEU:HD13	2.28	0.48
1:A:295:ASP:OD1	1:A:295:ASP:N	2.45	0.48
1:A:203:LEU:HD22	1:A:285:ILE:HD13	1.95	0.48
1:A:215:PHE:CD2	1:A:307:GLU:HG2	2.49	0.48
1:B:310:LEU:HD23	1:B:378:LEU:HD12	1.95	0.48
1:B:306:LEU:O	1:B:309:PRO:HD2	2.14	0.47
1:B:367:GLN:O	1:B:371:LEU:HD12	2.13	0.47
1:A:310:LEU:HD23	1:A:378:LEU:HD12	1.95	0.47
1:A:306:LEU:O	1:A:309:PRO:HD2	2.14	0.47
1:B:215:PHE:CD2	1:B:307:GLU:HG2	2.49	0.47
1:B:153:PHE:HB3	1:B:157:GLU:HB2	1.96	0.47
1:A:7:ILE:HG12	1:A:7:ILE:H	1.37	0.47
1:A:153:PHE:HB3	1:A:157:GLU:HB2	1.96	0.47
1:A:207:LEU:HD23	1:A:207:LEU:HA	1.77	0.46
1:B:210:PHE:CE2	1:B:378:LEU:HD11	2.42	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:ILE:HA	1:B:180:SER:OG	2.16	0.46
1:B:363:VAL:HG22	1:B:368:MET:HG2	1.98	0.46
1:B:176:CYS:O	1:B:180:SER:OG	2.27	0.46
1:A:339:ILE:HD12	1:A:339:ILE:H	1.80	0.46
1:B:150:LEU:HD23	1:B:150:LEU:HA	1.77	0.46
1:A:363:VAL:HG22	1:A:368:MET:HG2	1.98	0.46
1:B:339:ILE:H	1:B:339:ILE:HD12	1.80	0.46
1:A:404:LYS:HB3	1:A:404:LYS:HE2	1.63	0.45
1:B:7:ILE:HG12	1:B:7:ILE:H	1.37	0.45
1:A:150:LEU:CD1	1:B:36:SER:HB2	2.47	0.45
1:A:364:GLU:HG2	1:A:365:PRO:HD2	1.99	0.44
1:B:228:LYS:O	1:B:232:ASN:ND2	2.25	0.44
1:B:364:GLU:HG2	1:B:365:PRO:HD2	1.99	0.44
1:B:404:LYS:HB3	1:B:404:LYS:HE2	1.63	0.44
1:B:60:MET:O	1:B:64:VAL:HG23	2.18	0.44
1:A:178:LEU:HA	1:A:178:LEU:HD23	1.79	0.43
1:A:60:MET:O	1:A:64:VAL:HG23	2.18	0.43
1:A:308:THR:OG1	1:A:309:PRO:HD3	2.18	0.43
1:A:243:LYS:HA	1:A:279:ARG:HH21	1.84	0.43
1:B:20:ARG:HE	1:B:20:ARG:HB2	1.64	0.43
1:B:308:THR:OG1	1:B:309:PRO:HD3	2.18	0.43
1:B:377:MET:O	1:B:381:ILE:HG13	2.19	0.43
1:A:110:LEU:HD23	1:A:110:LEU:HA	1.86	0.42
1:B:140:ILE:HD13	1:B:140:ILE:HA	1.83	0.42
1:B:178:LEU:HD23	1:B:178:LEU:HA	1.79	0.42
1:A:364:GLU:OE1	1:A:366:LEU:HB2	2.20	0.42
1:A:377:MET:O	1:A:381:ILE:HG13	2.19	0.42
1:A:39:TYR:HD1	1:A:39:TYR:HA	1.71	0.42
1:B:336:SER:HA	1:B:339:ILE:HD13	2.01	0.42
1:A:66:PHE:HA	1:A:73:THR:HG22	2.01	0.42
1:B:70:VAL:HG22	1:B:180:SER:O	2.20	0.42
1:B:207:LEU:HD23	1:B:207:LEU:HA	1.77	0.42
1:A:20:ARG:HE	1:A:20:ARG:HB2	1.64	0.41
1:B:66:PHE:HA	1:B:73:THR:HG22	2.01	0.41
1:B:364:GLU:OE1	1:B:366:LEU:HB2	2.20	0.41
1:B:331:MET:HB3	1:B:401:PHE:CE1	2.55	0.41
1:B:354:LEU:HD23	1:B:359:ILE:HG21	2.02	0.41
1:A:378:LEU:HA	1:A:378:LEU:HD23	1.78	0.41
1:B:243:LYS:HA	1:B:279:ARG:HH21	1.84	0.41
1:A:32:LEU:HD23	1:A:32:LEU:HA	1.79	0.41
1:A:331:MET:HB3	1:A:401:PHE:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:LEU:HD23	1:A:359:ILE:HG21	2.02	0.41
1:B:32:LEU:HA	1:B:32:LEU:HD23	1.79	0.41
1:A:336:SER:HA	1:A:339:ILE:HD13	2.01	0.41
1:B:138:ILE:O	1:B:142:ILE:HG13	2.20	0.41
1:B:264:PHE:CD2	1:B:265:ARG:HG2	2.56	0.41
1:A:150:LEU:HA	1:A:150:LEU:HD23	1.77	0.41
1:B:356:LYS:HE3	1:B:356:LYS:HB2	1.85	0.41
1:A:264:PHE:CD2	1:A:265:ARG:HG2	2.56	0.41
1:A:138:ILE:O	1:A:142:ILE:HG13	2.20	0.40
1:B:288:LEU:HA	1:B:288:LEU:HD23	1.89	0.40
1:A:233:LEU:HB2	1:A:234:GLU:OE2	2.22	0.40
1:A:88:GLY:HA2	1:A:148:ILE:HD12	2.04	0.40
1:A:143:LEU:HD13	1:B:60:MET:HA	2.04	0.40
1:A:115:PHE:HB2	1:B:33:VAL:CG2	2.51	0.40
1:B:125:ARG:HH12	1:B:139:LEU:HD13	1.87	0.40
1:B:278:LEU:HD23	1:B:278:LEU:HA	1.84	0.40
1:B:330:GLN:HB3	1:B:332:ILE:HG12	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	383/509 (75%)	368 (96%)	15 (4%)	0	100	100
1	B	383/509 (75%)	368 (96%)	15 (4%)	0	100	100
All	All	766/1018 (75%)	736 (96%)	30 (4%)	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	339/443 (76%)	322 (95%)	17 (5%)	22	49
1	B	339/443 (76%)	323 (95%)	16 (5%)	23	50
All	All	678/886 (76%)	645 (95%)	33 (5%)	24	49

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	7	ILE
1	A	8	VAL
1	A	10	GLU
1	A	14	VAL
1	A	73	THR
1	A	180	SER
1	A	225	VAL
1	A	226	VAL
1	A	244	SER
1	A	304	LYS
1	A	320	SER
1	A	363	VAL
1	A	369	ILE
1	A	371	LEU
1	A	380	ASP
1	A	404	LYS
1	B	3	LYS
1	B	7	ILE
1	B	8	VAL
1	B	10	GLU
1	B	14	VAL
1	B	73	THR
1	B	225	VAL
1	B	226	VAL
1	B	244	SER
1	B	304	LYS

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Mol	Chain	Res	Type
1	B	320	SER
1	B	363	VAL
1	B	369	ILE
1	B	371	LEU
1	B	380	ASP
1	B	404	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	HIS
1	A	259	HIS
1	A	266	HIS
1	A	289	ASN
1	A	330	GLN
1	B	193	HIS
1	B	259	HIS
1	B	266	HIS
1	B	289	ASN
1	B	330	GLN

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 1 ligands modelled in this entry, 1 is 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.

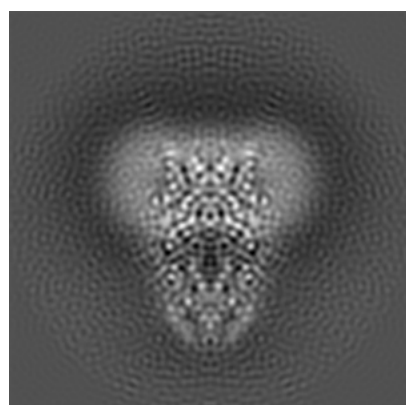
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32087. These allow visual inspection of the internal detail of the map and identification of artifacts.

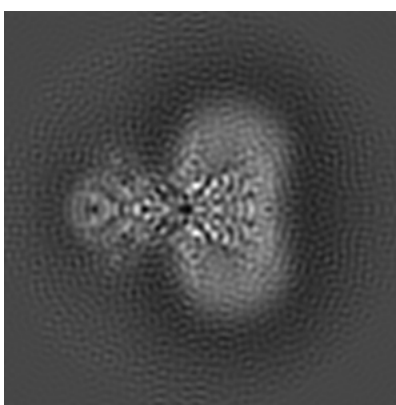
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

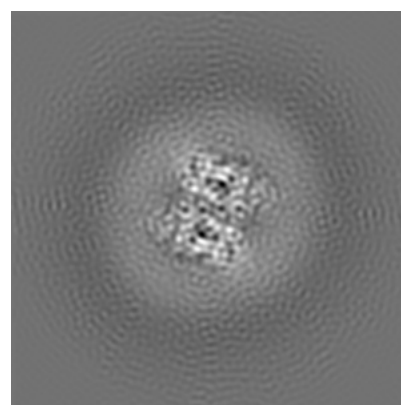
6.1.1 Primary map



X



Y

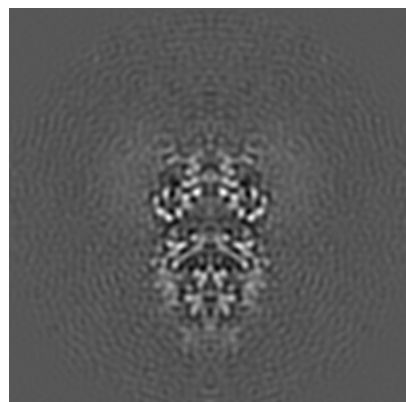


Z

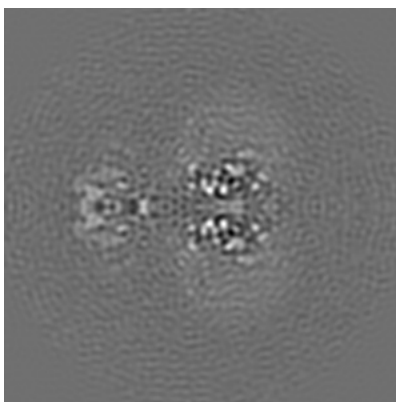
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

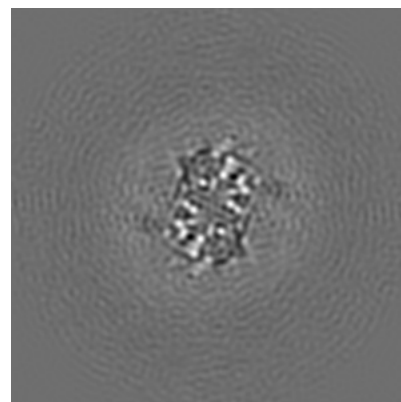
6.2.1 Primary map



X Index: 90



Y Index: 90

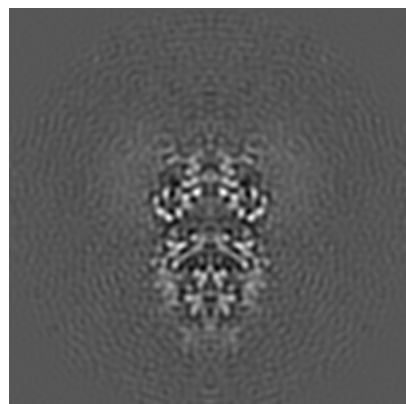


Z Index: 90

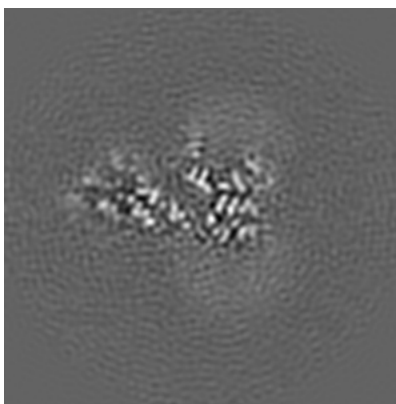
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

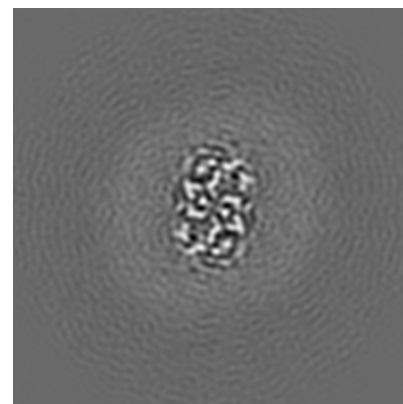
6.3.1 Primary map



X Index: 90



Y Index: 97

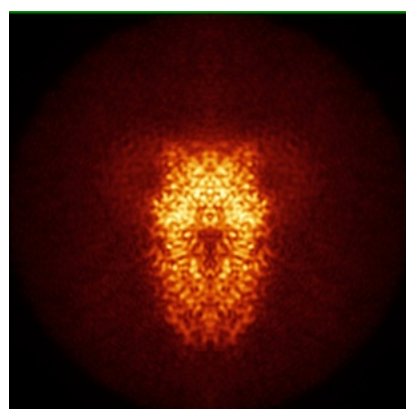


Z Index: 95

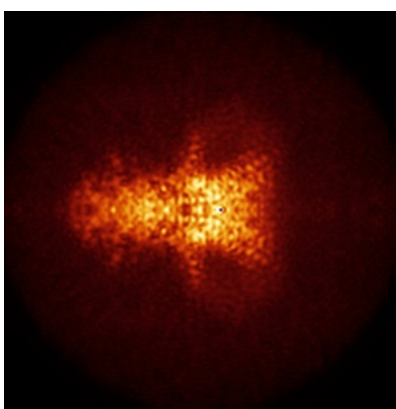
The images above show the largest variance slices of the map in three orthogonal directions.

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

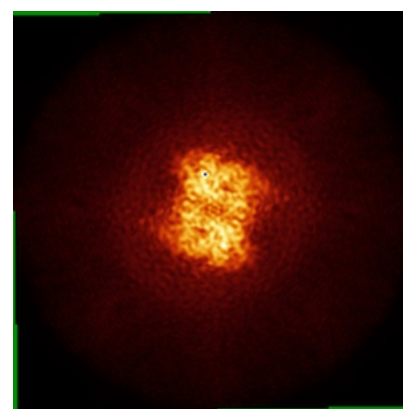
6.4.1 Primary map



X



Y

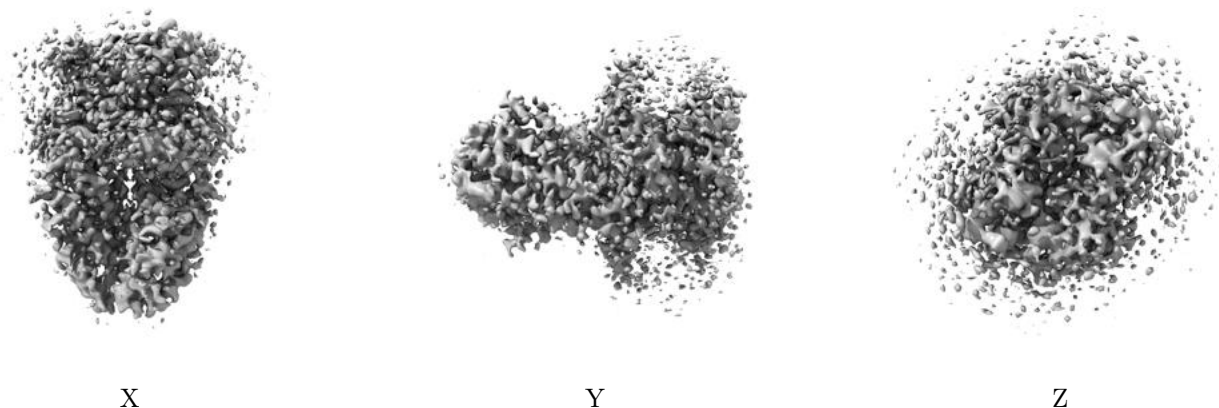


Z

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

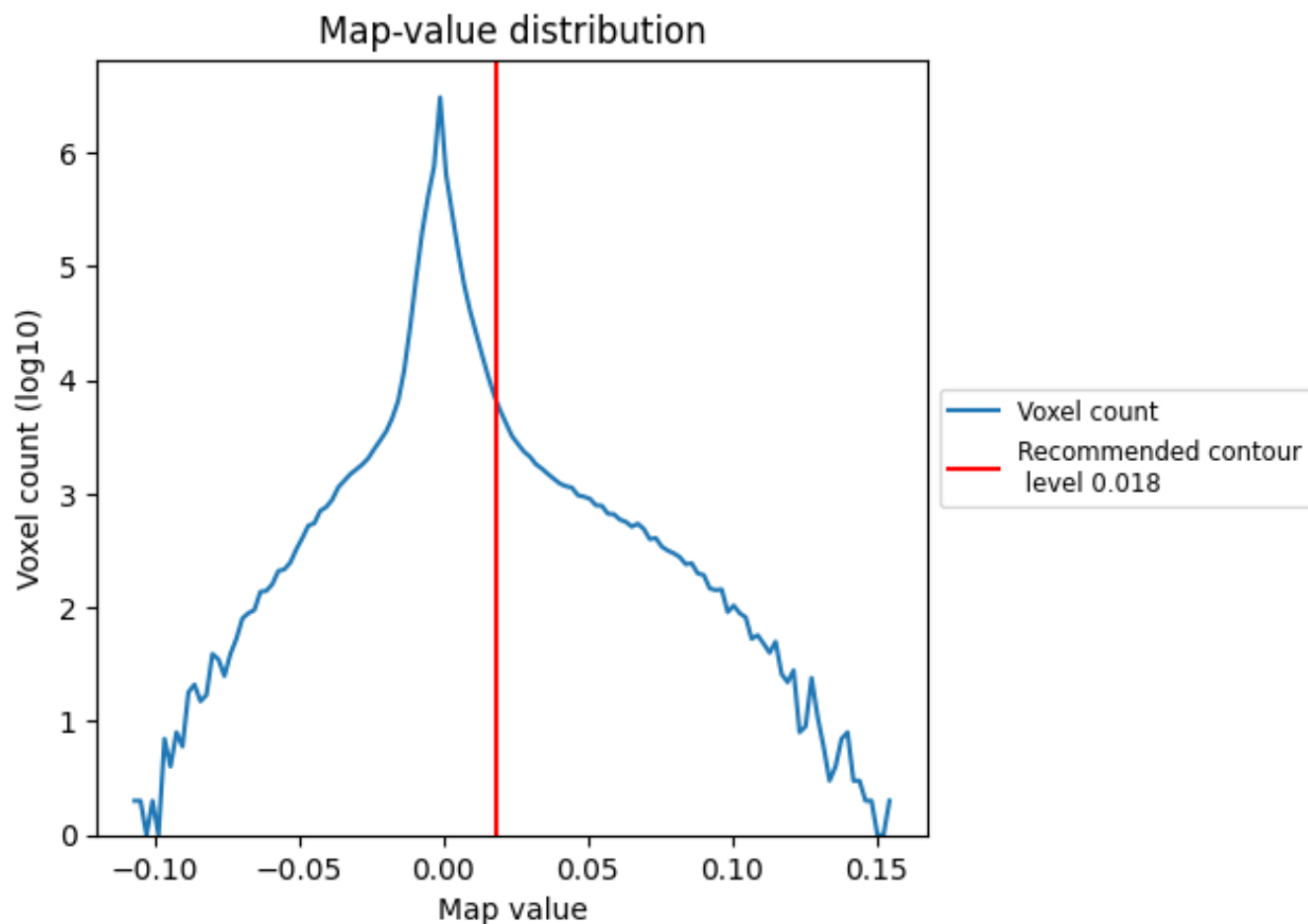
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

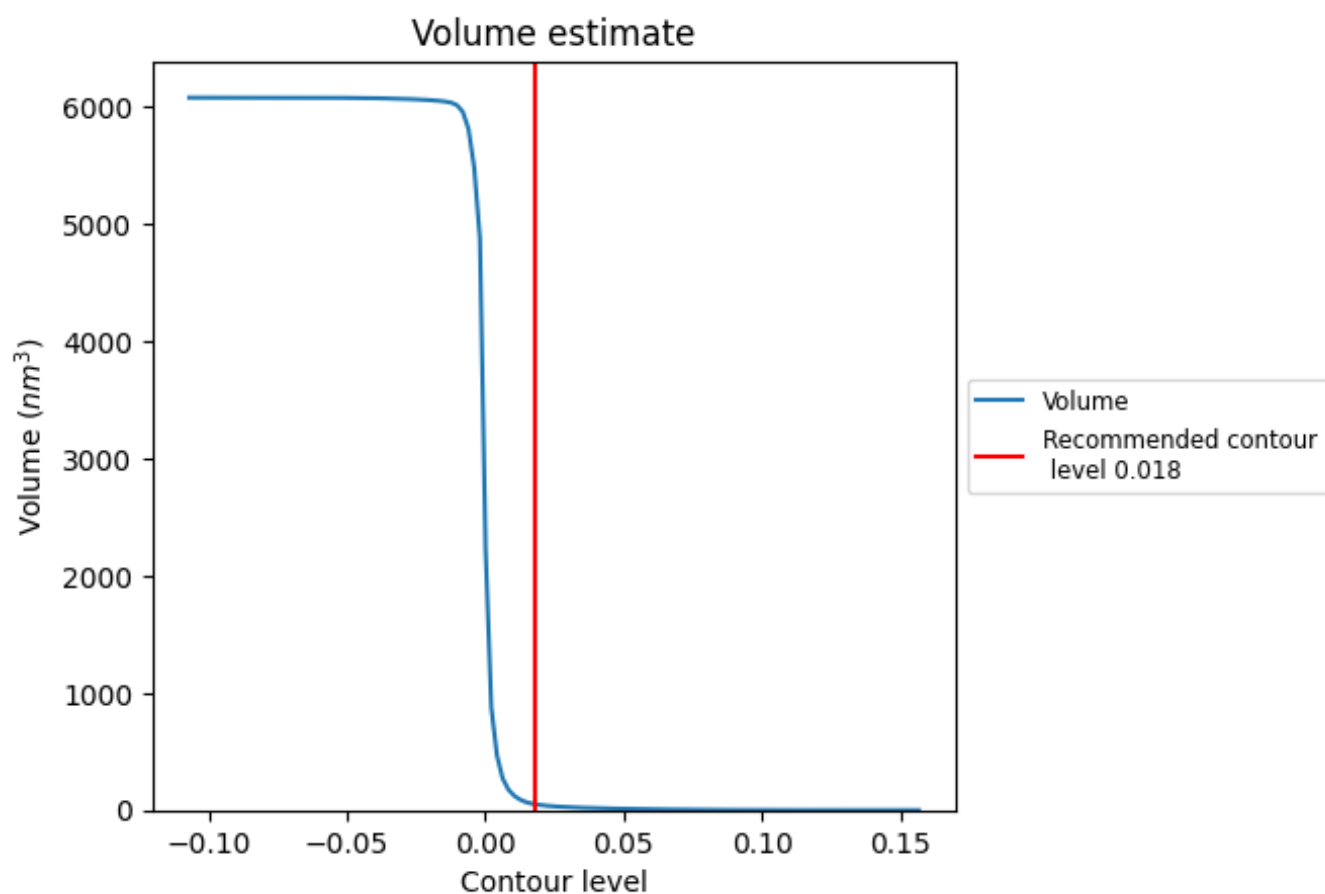
This section contains the results of statistical analysis of the map.

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.

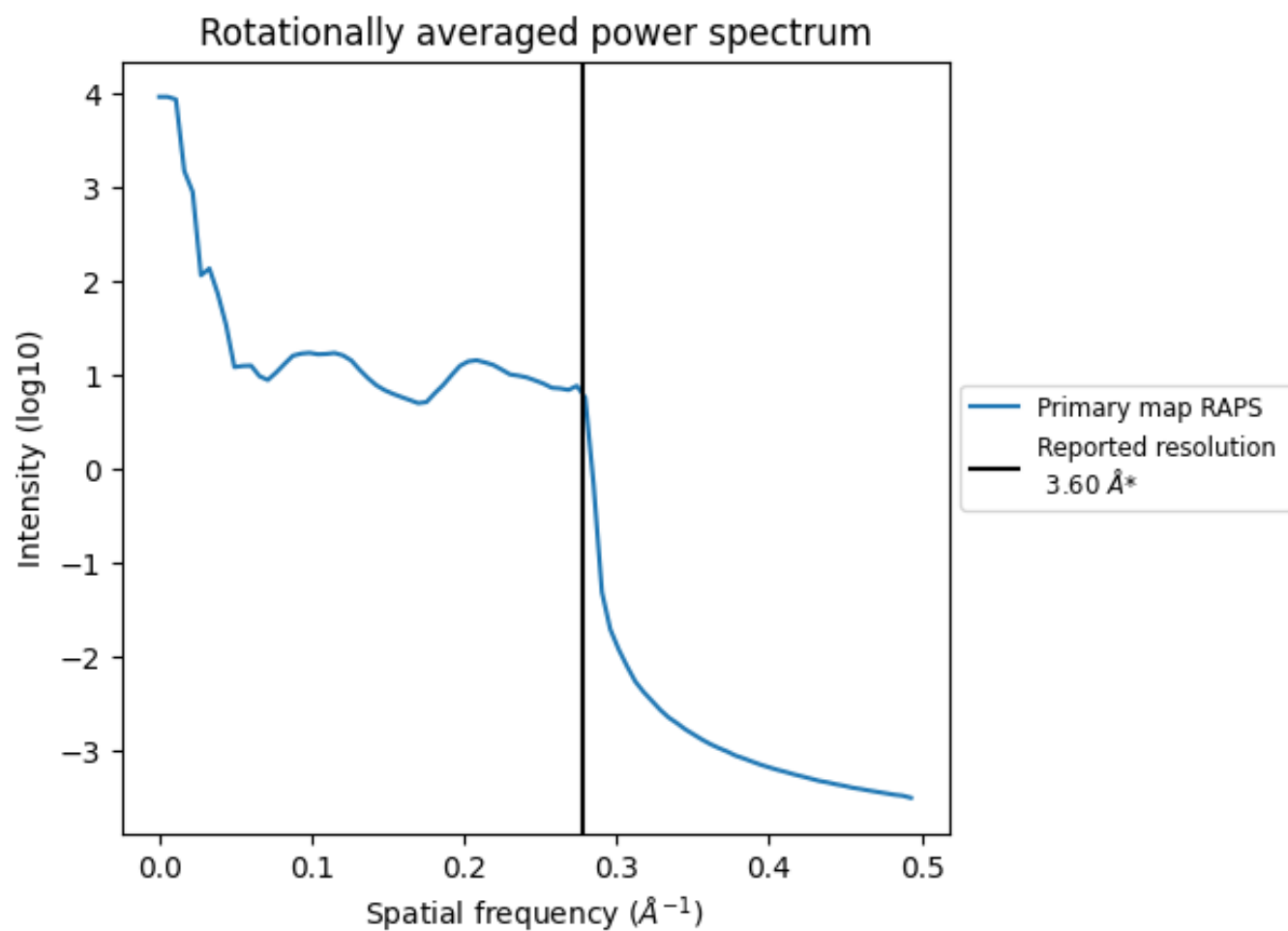
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 50 nm^3 ; this corresponds to an approximate mass of 45 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 0.278 Å⁻¹

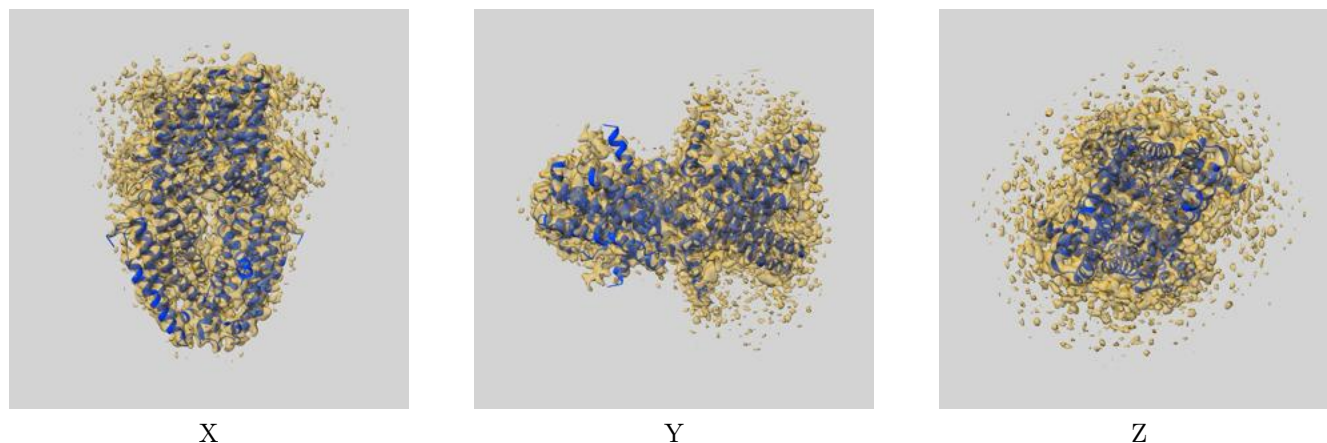
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

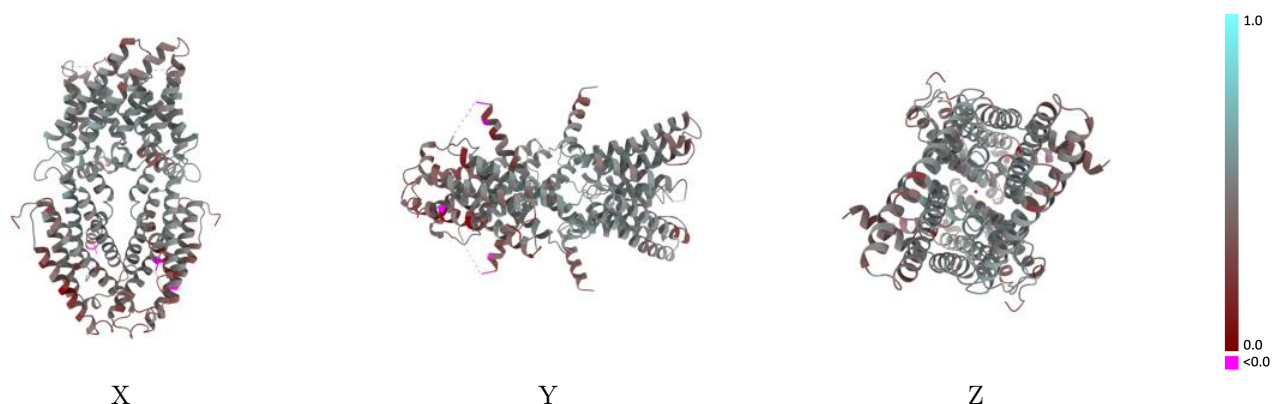
This section contains information regarding the fit between EMDB map EMD-32087 and PDB model 7VQ7. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

9.1 Map-model overlay [i](#)



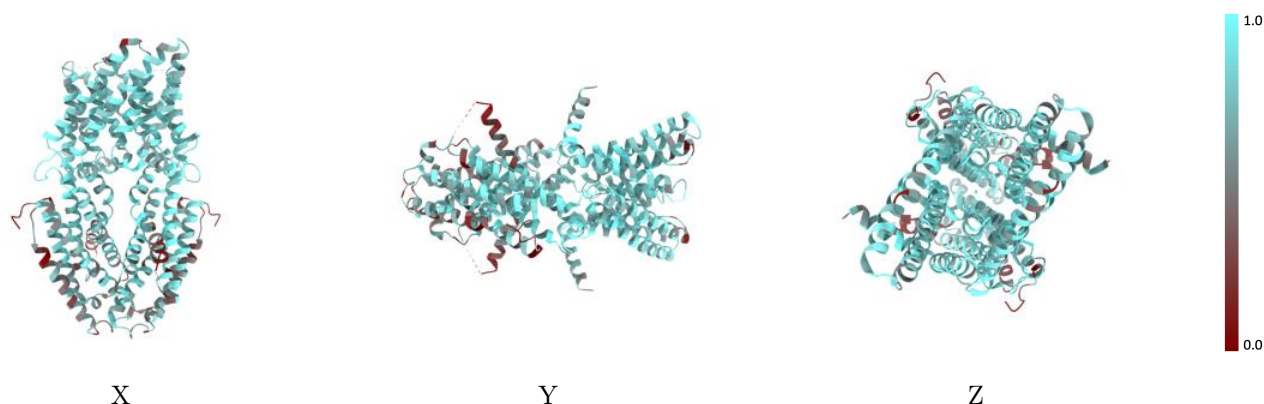
The images above show the 3D surface view of the map at the recommended contour level 0.018 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



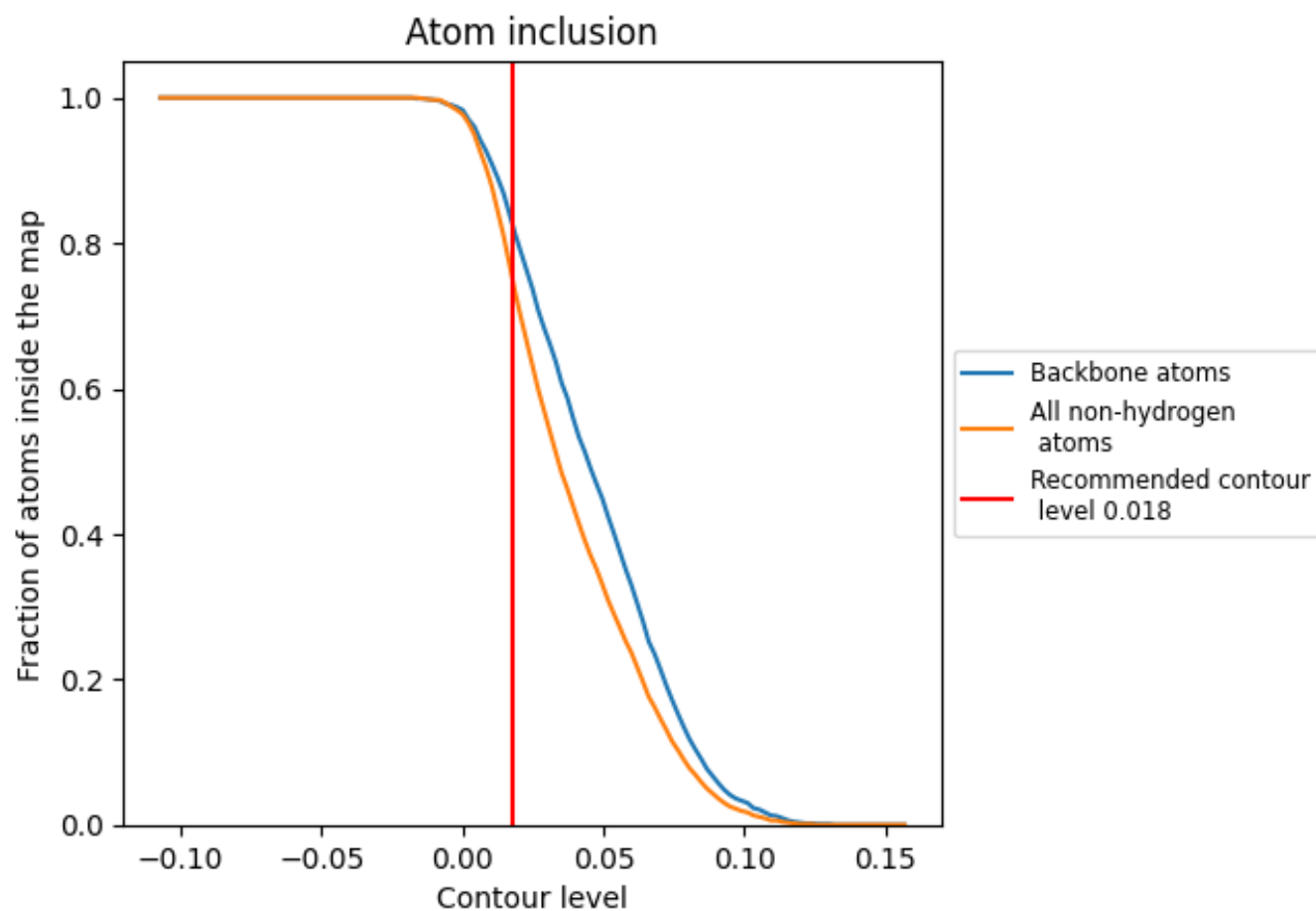
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.018).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 74% 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 (0.018) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7450	0.4480
A	0.7410	0.4470
B	0.7500	0.4490

