



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 10:48 PM UTC

PDB ID : 3J9S / pdb_00003j9s
EMDB ID : EMD-6272
Title : Single particle cryo-EM structure of rotavirus VP6 at 2.6 Angstrom resolution
Authors : Grant, T.; Grigorieff, N.
Deposited on : 2015-02-19
Resolution : 2.60 Å (reported)
Based on initial model : 1QHD

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
Mogul : 2022.3.0, CSD as543be (2022)
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

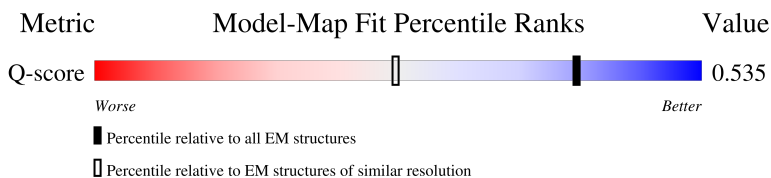
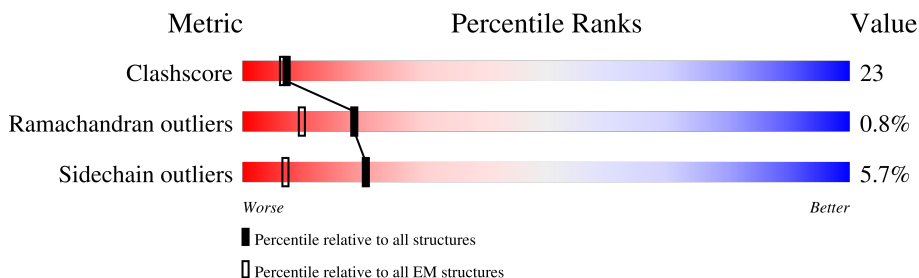
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.60 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	8728 (2.10 - 3.10)

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	398	11% 56% 39% ...

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3271 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 Intermediate capsid protein VP6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	398	Total	C	N	O	S	0	0
			3162	2004	546	597	15		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ACE	-	acetylation	UNP P04509

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
4	A	2	Total	Ca	0
			2	2	

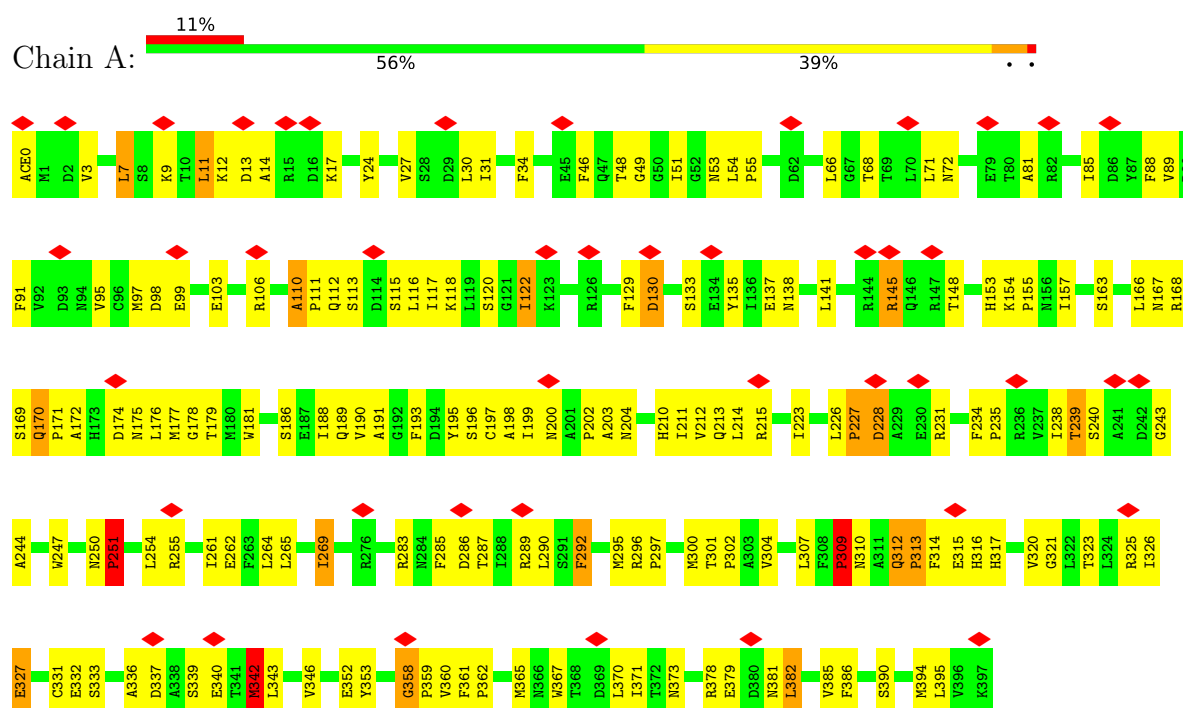
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		AltConf
5	A	105	Total	O	0
			105	105	

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: Intermediate capsid protein VP6



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	4000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	each particle	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	100	Depositor
Minimum defocus (nm)	400	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	29000	Depositor
Image detector	GATAN K2 (4k x 4k)	Depositor
Maximum map value	0.049	Depositor
Minimum map value	-0.020	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.0198	Depositor
Map size ($\text{\AA}$)	110.484, 110.484, 110.484	wwPDB
Map dimensions	108, 108, 108	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.023, 1.023, 1.023	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: CA, CL, ACE, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.42	19/3230 (0.6%)	1.20	34/4396 (0.8%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	110	ALA	C-N	6.46	1.41	1.33
1	A	361	PHE	C-N	6.18	1.41	1.33
1	A	234	PHE	C-N	6.08	1.40	1.33
1	A	13	ASP	C-O	-6.03	1.17	1.24
1	A	145	ARG	C-O	-5.89	1.16	1.23
1	A	12	LYS	C-O	-5.87	1.17	1.24
1	A	313	PRO	N-CD	5.79	1.55	1.47
1	A	7	LEU	C-O	-5.77	1.17	1.24
1	A	226	LEU	C-N	5.65	1.41	1.33
1	A	11	LEU	C-O	-5.39	1.17	1.24
1	A	97	MET	C-O	-5.39	1.17	1.24
1	A	331	CYS	C-O	-5.38	1.17	1.23
1	A	362	PRO	N-CD	5.32	1.55	1.47
1	A	359	PRO	N-CD	5.28	1.55	1.47
1	A	227	PRO	N-CD	5.25	1.55	1.47
1	A	170	GLN	C-N	5.17	1.41	1.34
1	A	171	PRO	N-CD	5.16	1.54	1.47
1	A	358	GLY	C-N	5.08	1.40	1.34
1	A	235	PRO	N-CD	5.07	1.54	1.47

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	361	PHE	CA-C-N	-8.93	113.23	119.66
1	A	361	PHE	C-N-CA	-8.93	113.23	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	339	SER	N-CA-C	8.37	120.49	111.36
1	A	310	ASN	N-CA-C	8.04	120.95	111.71
1	A	312	GLN	C-N-CD	7.90	137.97	120.60
1	A	342	MET	N-CA-C	7.88	119.95	111.36
1	A	106	ARG	N-CA-C	7.70	119.46	111.14
1	A	239	THR	N-CA-C	-7.07	101.71	110.41
1	A	244	ALA	N-CA-C	6.93	118.63	111.14
1	A	226	LEU	CA-C-N	-6.77	112.99	119.76
1	A	226	LEU	C-N-CA	-6.77	112.99	119.76
1	A	130	ASP	N-CA-C	6.71	120.65	111.39
1	A	110	ALA	CA-C-N	-6.64	112.92	119.90
1	A	110	ALA	C-N-CA	-6.64	112.92	119.90
1	A	122	ILE	N-CA-C	6.52	117.27	110.62
1	A	251	PRO	CA-N-CD	-6.38	103.07	112.00
1	A	153	HIS	N-CA-C	-6.36	98.88	109.24
1	A	309	PRO	CA-N-CD	-6.04	103.55	112.00
1	A	352	GLU	N-CA-C	5.99	118.59	111.71
1	A	135	TYR	N-CA-C	5.81	117.42	111.14
1	A	71	LEU	N-CA-C	5.79	117.67	111.36
1	A	17	LYS	N-CA-C	5.78	118.80	111.69
1	A	315	GLU	N-CA-C	5.72	117.60	111.36
1	A	283	ARG	N-CA-C	5.66	117.25	111.14
1	A	167	ASN	N-CA-C	5.60	117.46	111.36
1	A	170	GLN	CA-C-N	-5.54	112.90	119.05
1	A	170	GLN	C-N-CA	-5.54	112.90	119.05
1	A	234	PHE	CA-C-N	-5.47	114.16	119.90
1	A	234	PHE	C-N-CA	-5.47	114.16	119.90
1	A	358	GLY	CA-C-N	-5.41	113.05	119.05
1	A	358	GLY	C-N-CA	-5.41	113.05	119.05
1	A	269	ILE	N-CA-C	-5.16	102.85	109.30
1	A	72	ASN	N-CA-C	-5.10	101.45	109.50
1	A	327	GLU	N-CA-C	5.08	116.63	111.14

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	3162	0	3103	147	0
2	A	1	0	0	0	0
3	A	1	0	0	1	0
4	A	2	0	0	0	0
5	A	105	0	0	4	0
All	All	3271	0	3103	147	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 23.

All (147) 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:261:ILE:CD1	1:A:290:LEU:HD23	1.54	1.35
1:A:261:ILE:HD12	1:A:290:LEU:HD23	1.23	1.17
1:A:111:PRO:HB3	1:A:116:LEU:HD23	1.38	1.05
1:A:103:GLU:HG3	1:A:360:VAL:HG23	1.47	0.93
1:A:199:ILE:HB	1:A:296:ARG:NH1	1.84	0.92
1:A:175:ASN:ND2	1:A:314:PHE:HE2	1.68	0.92
1:A:261:ILE:HD12	1:A:290:LEU:CD2	2.01	0.91
1:A:170:GLN:OE1	1:A:175:ASN:HB3	1.69	0.91
1:A:213:GLN:HG2	5:A:645:HOH:O	1.70	0.90
1:A:163:SER:HB3	1:A:181:TRP:CZ2	2.08	0.88
1:A:103:GLU:HG3	1:A:360:VAL:CG2	2.03	0.88
1:A:262:GLU:OE1	1:A:269:ILE:HG12	1.74	0.87
1:A:138:ASN:OD1	1:A:148:THR:HB	1.75	0.86
1:A:261:ILE:CD1	1:A:290:LEU:CD2	2.49	0.85
1:A:169:SER:HA	1:A:176:LEU:CD2	2.06	0.84
1:A:99:GLU:OE1	1:A:113:SER:HB3	1.78	0.83
1:A:27:VAL:O	1:A:31:ILE:HG12	1.81	0.81
1:A:188:ILE:HD13	1:A:212:VAL:HG21	1.63	0.80
1:A:261:ILE:HD11	1:A:290:LEU:HD23	1.62	0.79
1:A:175:ASN:HD22	1:A:314:PHE:HE2	1.31	0.79
1:A:175:ASN:ND2	1:A:314:PHE:CE2	2.50	0.78
1:A:172:ALA:HB3	1:A:174:ASP:OD1	1.84	0.78
1:A:170:GLN:OE1	1:A:175:ASN:CB	2.32	0.77
1:A:148:THR:O	1:A:332:GLU:HG2	1.85	0.76
1:A:227:PRO:O	1:A:228:ASP:HB2	1.87	0.75
1:A:337:ASP:OD1	1:A:340:GLU:HB2	1.88	0.73
1:A:170:GLN:OE1	1:A:175:ASN:N	2.22	0.72
1:A:307:LEU:HD13	1:A:316:HIS:ND1	2.05	0.72
1:A:169:SER:HA	1:A:176:LEU:HD21	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:MET:CE	1:A:386:PHE:O	2.38	0.72
1:A:103:GLU:OE2	1:A:358:GLY:HA3	1.91	0.70
1:A:46:PHE:CE2	1:A:118:LYS:HD3	2.26	0.69
1:A:166:LEU:HD22	1:A:176:LEU:HD22	1.74	0.69
1:A:170:GLN:OE1	1:A:175:ASN:CA	2.41	0.69
1:A:172:ALA:CB	1:A:174:ASP:OD1	2.41	0.68
1:A:297:PRO:HD2	1:A:300:MET:HE2	1.75	0.67
1:A:265:LEU:HA	1:A:286:ASP:OD1	1.94	0.67
1:A:168:ARG:O	1:A:176:LEU:HD23	1.94	0.67
1:A:190:VAL:HG21	1:A:210:HIS:HB2	1.78	0.66
1:A:178:GLY:O	1:A:193:PHE:CD2	2.49	0.65
1:A:312:GLN:HB3	1:A:313:PRO:HA	1.78	0.65
1:A:9:LYS:HD3	1:A:394:MET:HE1	1.78	0.65
1:A:48:THR:HG22	1:A:115:SER:OG	1.96	0.65
1:A:238:ILE:HG22	1:A:239:THR:O	1.96	0.65
1:A:99:GLU:OE1	1:A:113:SER:CB	2.44	0.65
1:A:289:ARG:HG3	1:A:289:ARG:O	1.96	0.65
1:A:157:ILE:HG13	1:A:157:ILE:O	1.97	0.64
1:A:228:ASP:N	1:A:321:GLY:O	2.28	0.64
1:A:113:SER:O	1:A:117:ILE:HG13	1.97	0.64
1:A:178:GLY:O	1:A:193:PHE:CE2	2.52	0.63
1:A:169:SER:HA	1:A:176:LEU:HD23	1.78	0.63
1:A:91:PHE:O	1:A:95:VAL:HG23	1.98	0.63
1:A:250:ASN:O	1:A:317:HIS:ND1	2.33	0.61
1:A:133:SER:O	1:A:137:GLU:HG3	2.01	0.61
1:A:154:LYS:HE3	3:A:502:CL:CL	2.38	0.61
1:A:254:LEU:O	1:A:320:VAL:HG12	2.02	0.60
1:A:269:ILE:H	1:A:269:ILE:HD12	1.66	0.60
1:A:170:GLN:NE2	1:A:175:ASN:O	2.35	0.59
1:A:373:ASN:O	1:A:378:ARG:NH1	2.36	0.59
1:A:9:LYS:HD3	1:A:394:MET:CE	2.32	0.59
1:A:342:MET:HE1	1:A:386:PHE:O	2.01	0.59
1:A:81:ALA:O	1:A:85:ILE:HG13	2.03	0.58
1:A:168:ARG:O	1:A:176:LEU:CD2	2.52	0.57
1:A:261:ILE:HD13	1:A:290:LEU:HD23	1.71	0.57
1:A:227:PRO:O	1:A:228:ASP:CB	2.52	0.57
1:A:3:VAL:HG21	5:A:620:HOH:O	2.06	0.56
1:A:196:SER:O	1:A:197:CYS:HB3	2.04	0.56
1:A:202:PRO:O	1:A:203:ALA:HB3	2.06	0.56
1:A:214:LEU:N	1:A:214:LEU:HD23	2.21	0.56
1:A:262:GLU:OE1	1:A:269:ILE:CG1	2.52	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:TYR:OH	1:A:68:THR:HG22	2.07	0.55
1:A:255:ARG:NH1	1:A:255:ARG:HG3	2.22	0.55
1:A:292:PHE:CD1	1:A:292:PHE:N	2.75	0.55
1:A:163:SER:HB3	1:A:181:TRP:CE2	2.42	0.54
1:A:312:GLN:HG2	1:A:314:PHE:CE2	2.42	0.54
1:A:312:GLN:CB	1:A:313:PRO:HA	2.37	0.53
1:A:197:CYS:O	1:A:198:ALA:HB3	2.09	0.53
1:A:360:VAL:HG11	1:A:381:ASN:CG	2.33	0.53
1:A:122:ILE:HD12	1:A:122:ILE:O	2.09	0.52
1:A:247:TRP:H	1:A:247:TRP:CD1	2.27	0.52
1:A:239:THR:CG2	1:A:240:SER:N	2.73	0.52
1:A:239:THR:HG22	1:A:240:SER:N	2.24	0.52
1:A:367:TRP:O	1:A:371:ILE:HG22	2.09	0.52
1:A:14:ALA:HA	1:A:30:LEU:HD21	1.93	0.51
1:A:200:ASN:HB2	5:A:696:HOH:O	2.10	0.51
1:A:177:MET:HA	1:A:193:PHE:O	2.11	0.51
1:A:342:MET:HE1	1:A:390:SER:HB3	1.93	0.50
1:A:0:ACE:CH3	1:A:120:SER:HB3	2.41	0.50
1:A:307:LEU:HD13	1:A:316:HIS:CG	2.46	0.50
1:A:231:ARG:HH11	1:A:231:ARG:HG3	1.77	0.50
1:A:34:PHE:CD2	1:A:66:LEU:HD11	2.47	0.50
1:A:110:ALA:O	1:A:112:GLN:NE2	2.45	0.49
1:A:145:ARG:O	1:A:145:ARG:HG3	2.12	0.49
1:A:51:ILE:O	1:A:51:ILE:HG23	2.11	0.49
1:A:211:ILE:HG12	1:A:289:ARG:HB2	1.95	0.49
1:A:301:THR:HB	1:A:302:PRO:CD	2.43	0.49
1:A:301:THR:HB	1:A:302:PRO:HD2	1.94	0.49
1:A:49:GLY:N	1:A:98:ASP:OD2	2.37	0.48
1:A:85:ILE:O	1:A:89:VAL:HG23	2.14	0.48
1:A:170:GLN:CD	1:A:175:ASN:O	2.56	0.48
1:A:223:ILE:HG23	1:A:325:ARG:O	2.14	0.48
1:A:24:TYR:OH	1:A:68:THR:CG2	2.62	0.48
1:A:53:ASN:HB2	1:A:353:TYR:O	2.14	0.47
1:A:129:PHE:CD2	1:A:130:ASP:HB2	2.50	0.47
1:A:382:LEU:O	1:A:385:VAL:HG22	2.15	0.47
1:A:46:PHE:CD2	1:A:118:LYS:HD3	2.49	0.47
1:A:175:ASN:HB2	1:A:312:GLN:NE2	2.29	0.47
1:A:103:GLU:HG3	1:A:360:VAL:CB	2.44	0.46
1:A:7:LEU:O	1:A:11:LEU:HG	2.14	0.46
1:A:155:PRO:O	1:A:186:SER:HB3	2.15	0.46
1:A:177:MET:O	1:A:177:MET:CG	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:PRO:O	1:A:309:PRO:HD2	2.16	0.46
1:A:327:GLU:CD	5:A:609:HOH:O	2.59	0.46
1:A:239:THR:CG2	1:A:240:SER:H	2.28	0.46
1:A:336:ALA:HB2	1:A:343:LEU:HD22	1.97	0.46
1:A:170:GLN:N	1:A:170:GLN:HE21	2.14	0.46
1:A:24:TYR:OH	1:A:68:THR:HB	2.16	0.46
1:A:255:ARG:HG3	1:A:255:ARG:HH11	1.81	0.45
1:A:333:SER:OG	1:A:379:GLU:OE2	2.35	0.45
1:A:170:GLN:OE1	1:A:175:ASN:C	2.60	0.45
1:A:264:LEU:O	1:A:286:ASP:OD1	2.35	0.45
1:A:255:ARG:HH11	1:A:255:ARG:CG	2.29	0.45
1:A:176:LEU:O	1:A:195:TYR:CD1	2.70	0.44
1:A:231:ARG:HG3	1:A:231:ARG:NH1	2.32	0.44
1:A:215:ARG:HG3	1:A:371:ILE:O	2.17	0.44
1:A:170:GLN:NE2	1:A:176:LEU:HD23	2.32	0.44
1:A:342:MET:HE3	1:A:342:MET:HB3	1.59	0.44
1:A:196:SER:O	1:A:197:CYS:CB	2.66	0.44
1:A:203:ALA:O	1:A:204:ASN:HB2	2.18	0.43
1:A:169:SER:C	1:A:170:GLN:HE21	2.26	0.43
1:A:214:LEU:HD11	1:A:285:PHE:CE2	2.54	0.42
1:A:307:LEU:CD1	1:A:316:HIS:CG	3.02	0.42
1:A:11:LEU:HD21	1:A:88:PHE:HB3	2.00	0.42
1:A:342:MET:H	1:A:342:MET:HG2	1.61	0.42
1:A:382:LEU:HD23	1:A:382:LEU:HA	1.93	0.42
1:A:141:LEU:HB2	1:A:148:THR:HG22	2.02	0.42
1:A:251:PRO:HA	1:A:317:HIS:O	2.20	0.41
1:A:189:GLN:HG3	1:A:323:THR:HG23	2.01	0.41
1:A:11:LEU:HD12	1:A:395:LEU:HD11	2.00	0.41
1:A:304:VAL:O	1:A:307:LEU:HB2	2.20	0.41
1:A:213:GLN:HG3	1:A:287:THR:HG23	2.02	0.41
1:A:54:LEU:HA	1:A:55:PRO:HD3	1.89	0.41
1:A:240:SER:O	1:A:243:GLY:N	2.54	0.41
1:A:197:CYS:O	1:A:197:CYS:SG	2.78	0.41
1:A:346:VAL:HG12	1:A:365:MET:HE2	2.03	0.41
1:A:54:LEU:HD12	1:A:55:PRO:HD2	2.03	0.40
1:A:179:THR:HA	1:A:191:ALA:O	2.21	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	396/398 (100%)	388 (98%)	5 (1%)	3 (1%)	16	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	309	PRO
1	A	228	ASP
1	A	251	PRO

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	106/350 (30%)	100 (94%)	6 (6%)	18	40

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	292	PHE
1	A	295	MET
1	A	326	ILE
1	A	342	MET
1	A	370	LEU
1	A	382	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 4 ligands modelled in this entry, 4 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 [i](#)

There are no chain breaks in this entry.

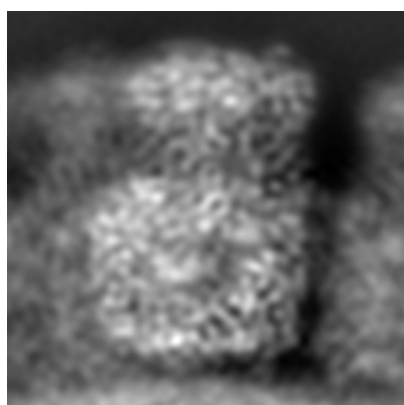
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6272. These allow visual inspection of the internal detail of the map and identification of artifacts.

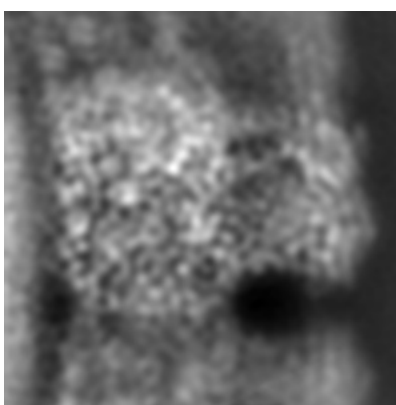
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

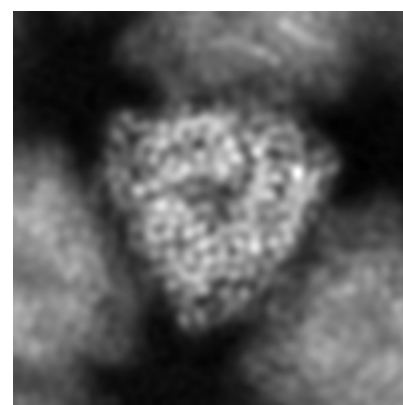
6.1.1 Primary map



X



Y

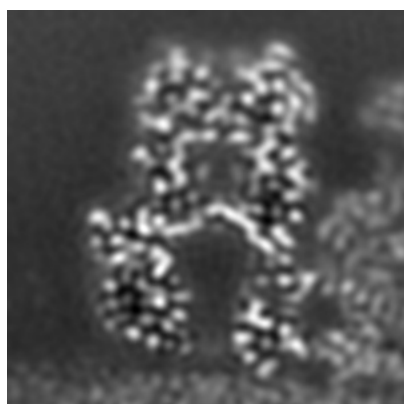


Z

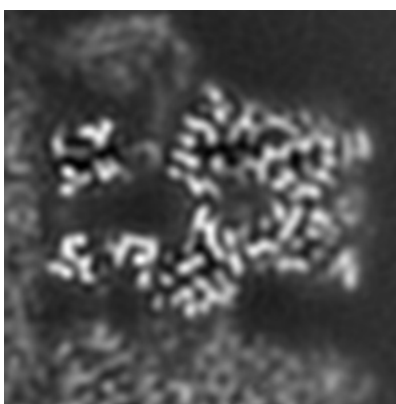
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

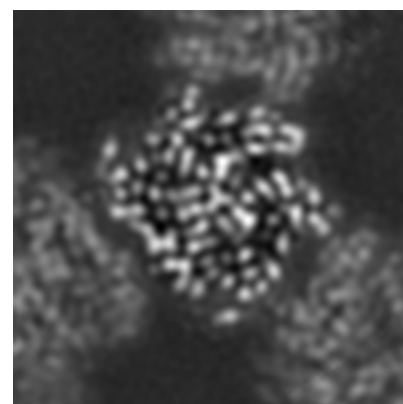
6.2.1 Primary map



X Index: 54



Y Index: 54

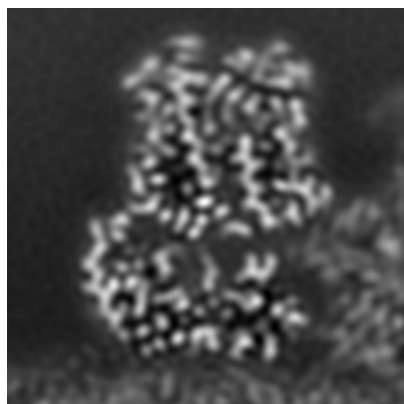


Z Index: 54

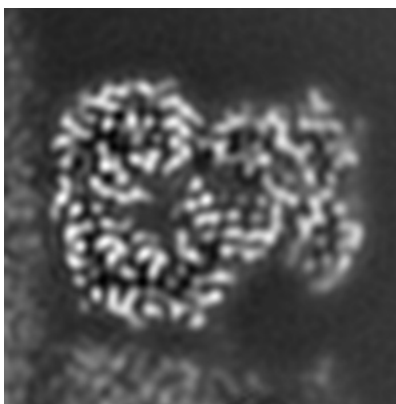
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

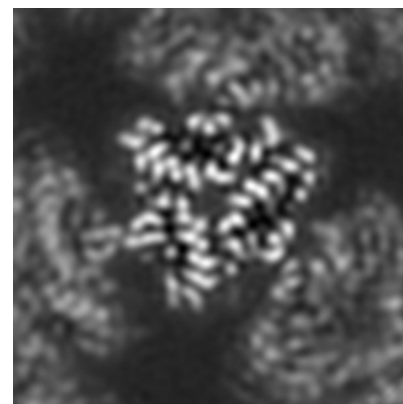
6.3.1 Primary map



X Index: 46



Y Index: 64

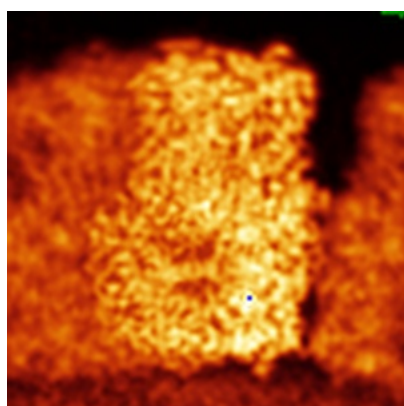


Z Index: 19

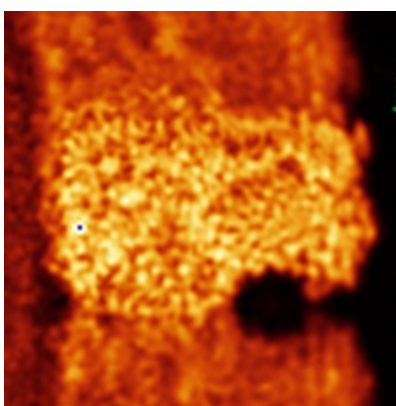
The images above show the largest variance slices of the map in three orthogonal directions.

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

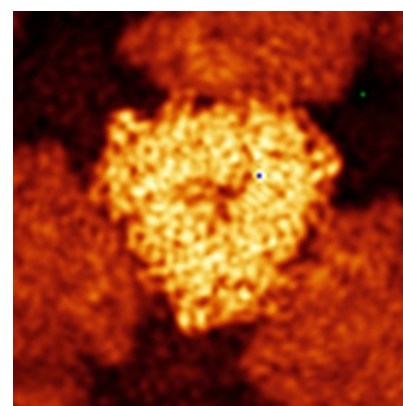
6.4.1 Primary map



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

This section was not generated.

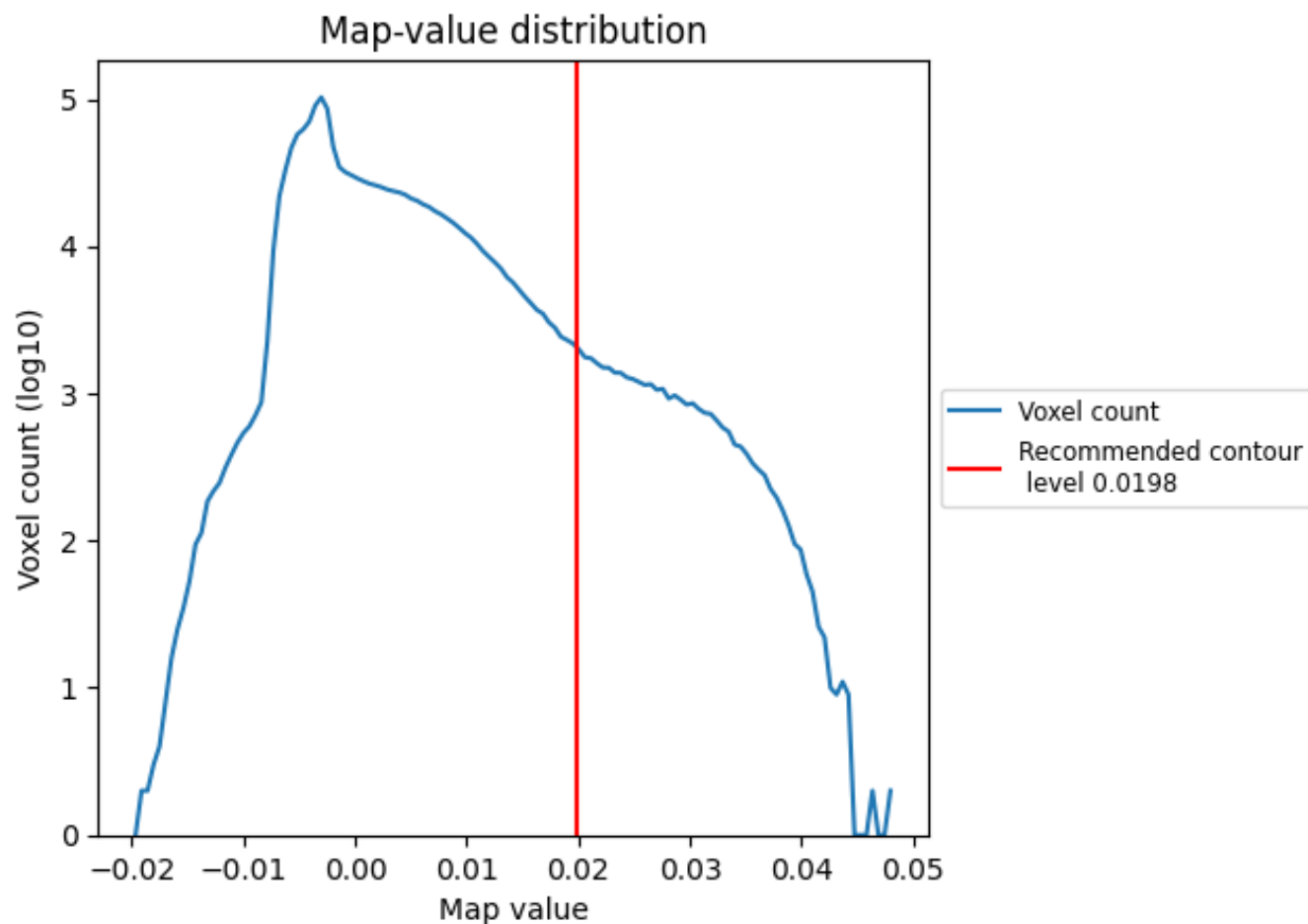
6.6 Mask visualisation

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

7 Map analysis [i](#)

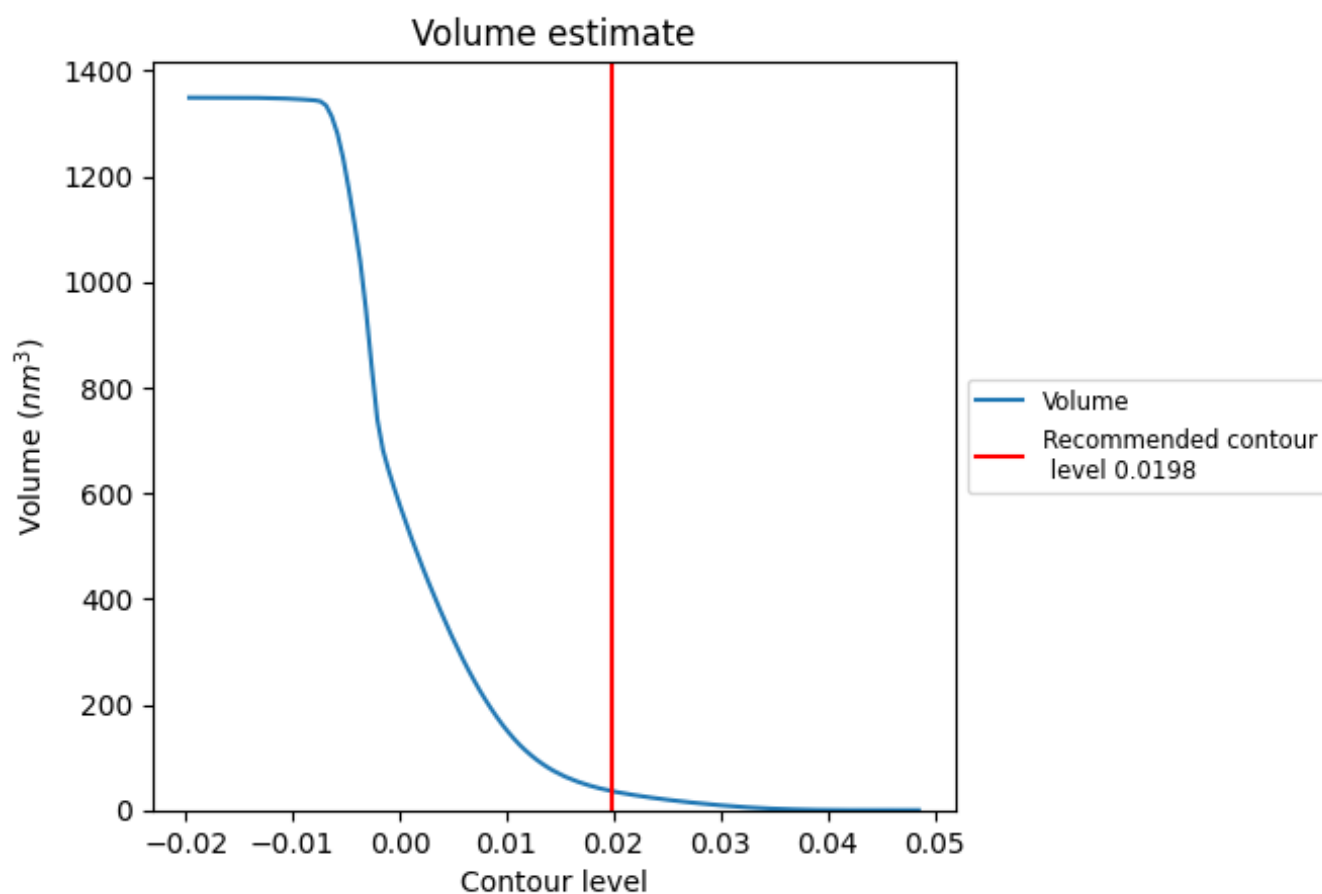
This section contains the results of statistical analysis of the map.

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.

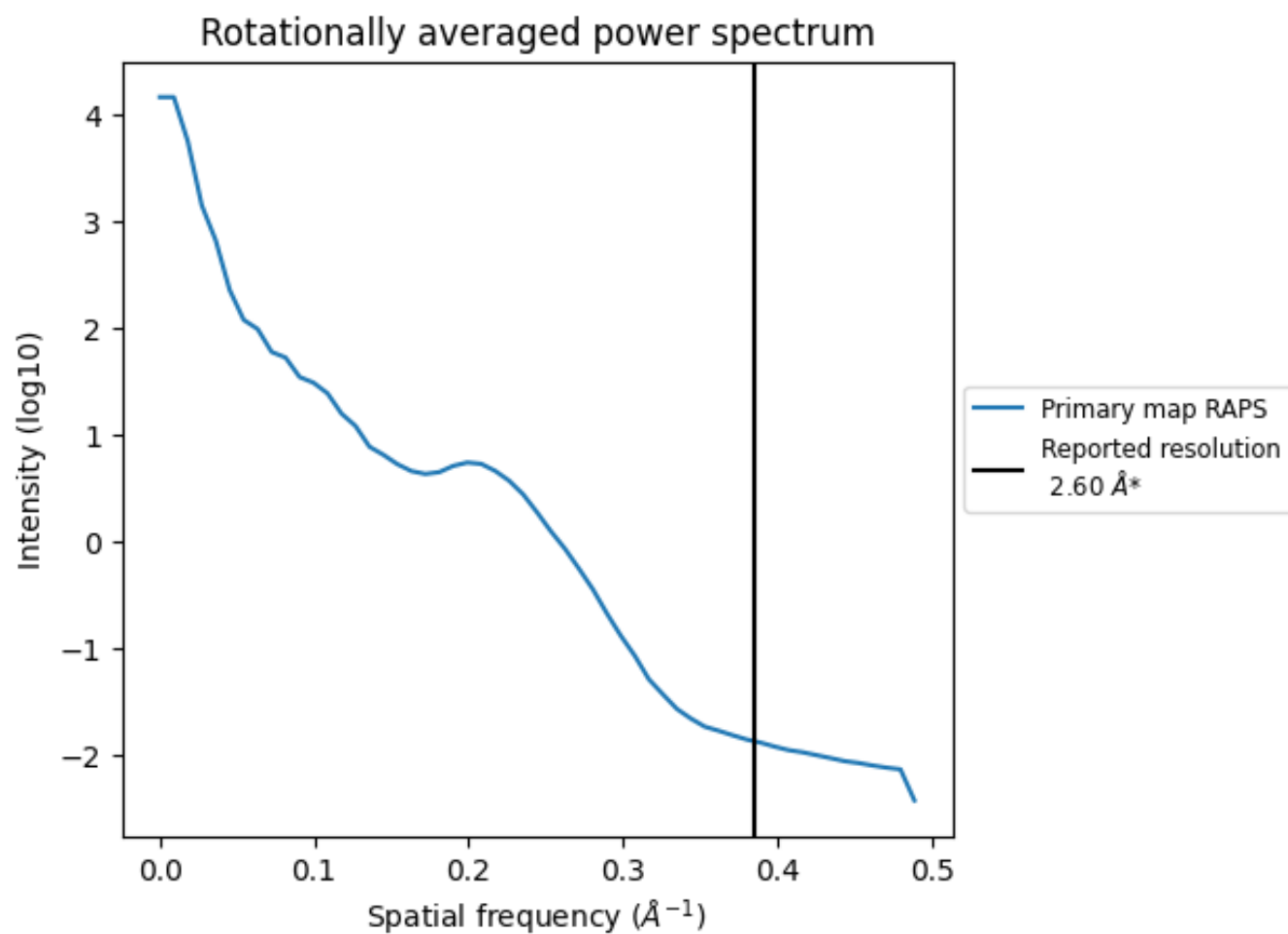
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 36 nm³; this corresponds to an approximate mass of 33 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 0.385 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

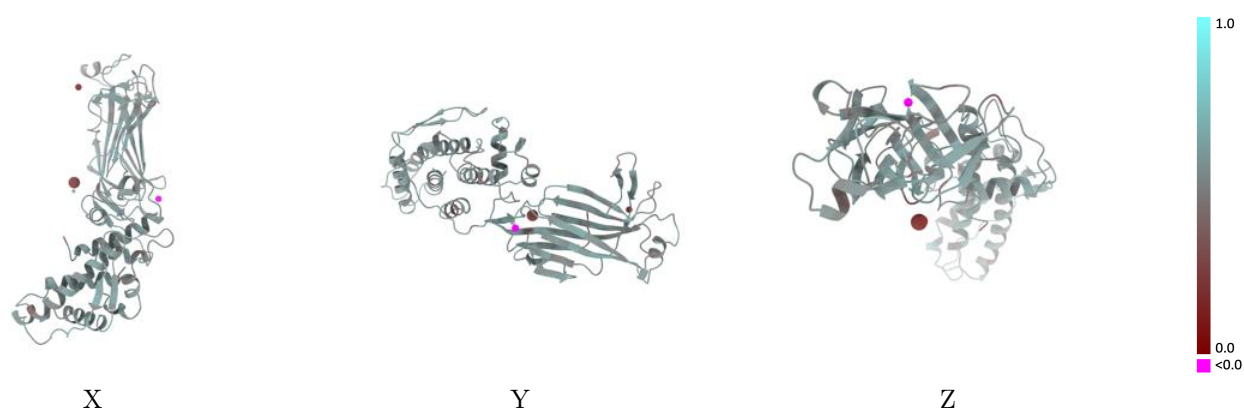
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-6272 and PDB model 3J9S. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)

This section was not generated.

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

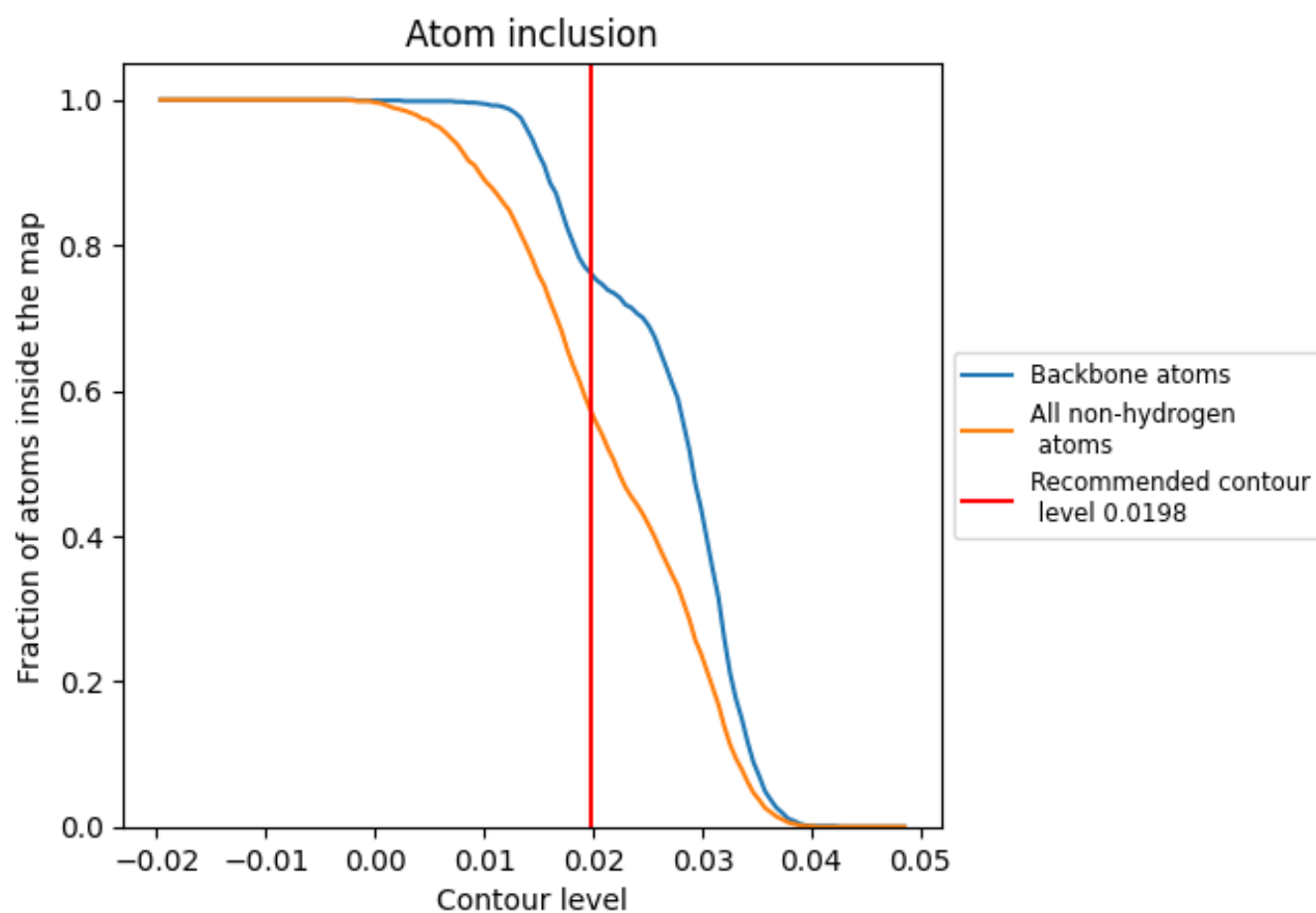


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 57% 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 (0.0198) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5710	0.5350
A	0.5710	0.5350

