



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 08:27 PM UTC

PDB ID : 6VQH / pdb_00006vqh
EMDB ID : EMD-21350
Title : Mammalian V-ATPase from rat brain membrane-embedded Vo region rotational state 3 (from focused refinement)
Authors : Abbas, Y.M.; Rubinstein, J.L.
Deposited on : 2020-02-05
Resolution : 4.40 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:

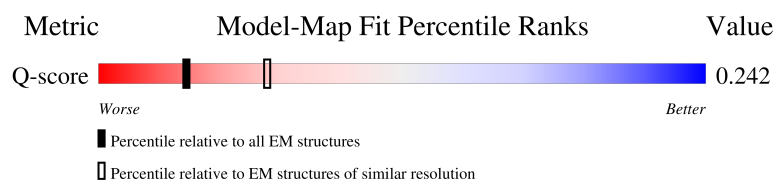
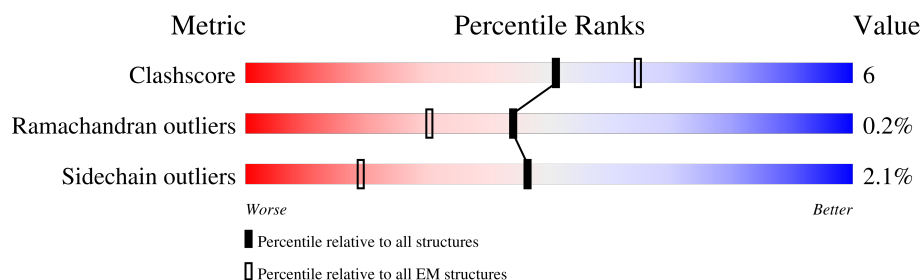
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 4.40 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	3132 (3.91 - 4.90)

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	H	247	
2	L	119	
3	a	838	
4	b	205	

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Mol	Chain	Length	Quality of chain
5	c	463	 8% 91%
6	d	351	 24% 87% 12%
7	e	81	 65% 94%
8	f	98	 64% 86% 14%
9	g	155	 21% 81% 13%
9	h	155	 19% 81% 14%
9	i	155	 23% 78% 15%
9	j	155	 21% 81% 14%
9	k	155	 25% 82% 14%
9	l	155	 22% 85% 12%
9	m	155	 25% 78% 16%
9	n	155	 21% 82% 14%
9	o	155	 30% 72% 24%
10	p	350	 13% 85%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 19338 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 ATPase H⁺-transporting V1 subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	H	104	Total	C	N	O	S	0	0
			803	509	138	153	3		

- Molecule 2 is a protein called V-type proton ATPase subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	L	110	Total	C	N	O	S	0	0
			875	553	157	163	2		

- Molecule 3 is a protein called V-type proton ATPase 116 kDa subunit a isoform 1.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	a	441	Total	C	N	O	0	0
			2173	1291	441	441		

- Molecule 4 is a protein called ATPase, H⁺ transporting, V0 subunit B (Predicted), isoform CRA_a.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	b	203	Total	C	N	O	S	0	0
			1503	996	237	259	11		

- Molecule 5 is a protein called V-type proton ATPase subunit S1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	c	41	Total	C	N	O	S	0	0
			337	228	51	54	4		

- Molecule 6 is a protein called V-type proton ATPase subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	d	350	Total	C	N	O	S	0	0
			2833	1829	460	530	14		

- Molecule 7 is a protein called V-type proton ATPase subunit e 2.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	e	78	Total	C	N	O	0	0
			384	228	78	78		

- Molecule 8 is a protein called Ribonuclease K.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	f	84	Total	C	N	O	0	0
			412	244	84	84		

- Molecule 9 is a protein called V-type proton ATPase 16 kDa proteolipid subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	g	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	h	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	i	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	j	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	k	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	l	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	m	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	n	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		
9	o	150	Total	C	N	O	S	0	0
			1068	699	171	190	8		

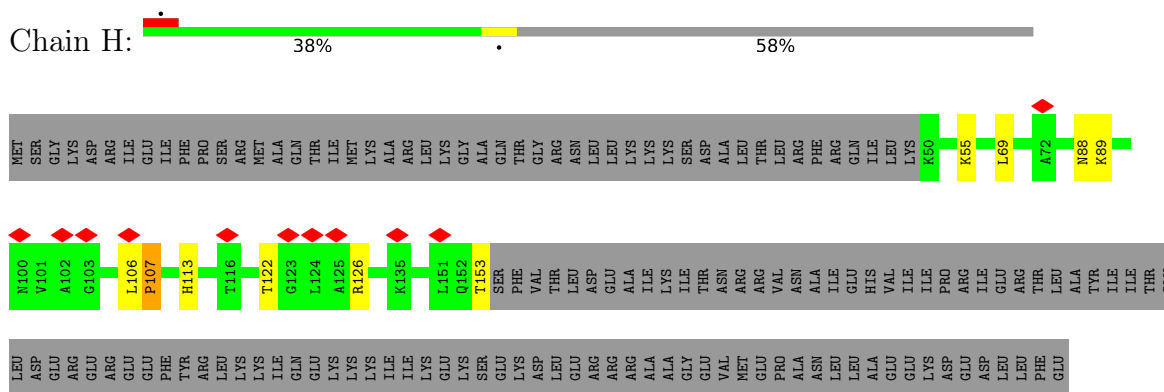
- Molecule 10 is a protein called Renin receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	p	52	Total	C	N	O	S	0	0
			406	269	61	73	3		

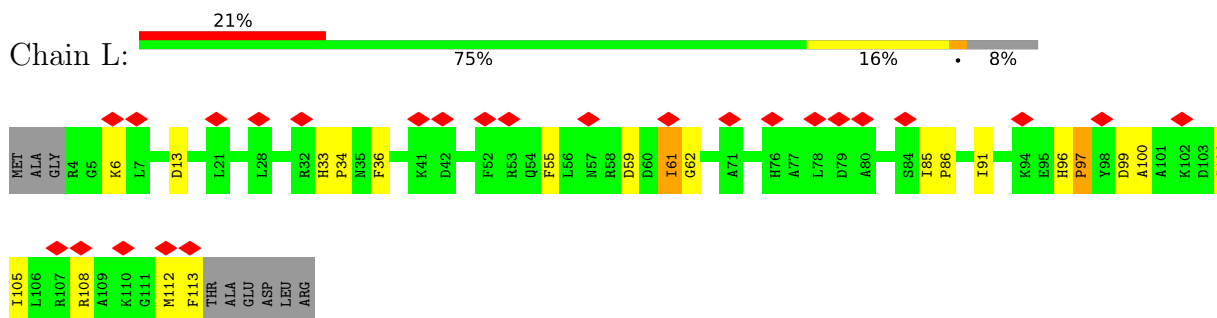
3 Residue-property plots

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.

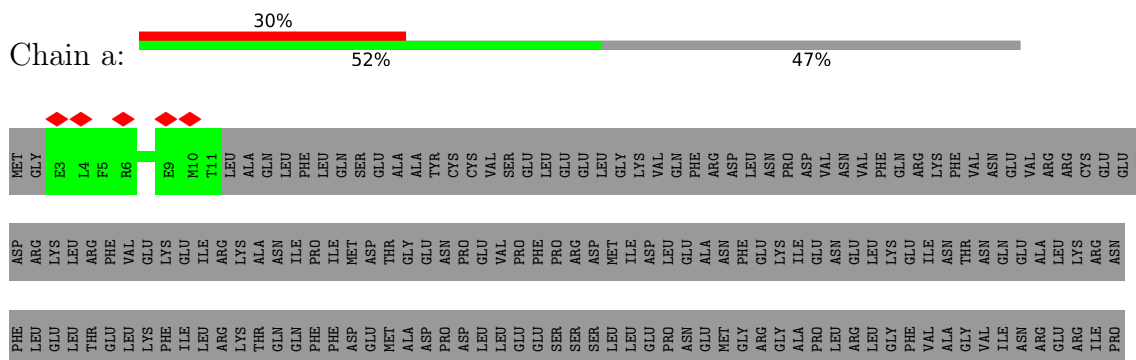
- Molecule 1: ATPase H⁺-transporting V1 subunit D

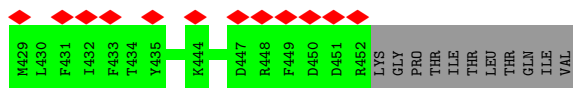


- Molecule 2: V-type proton ATPase subunit F

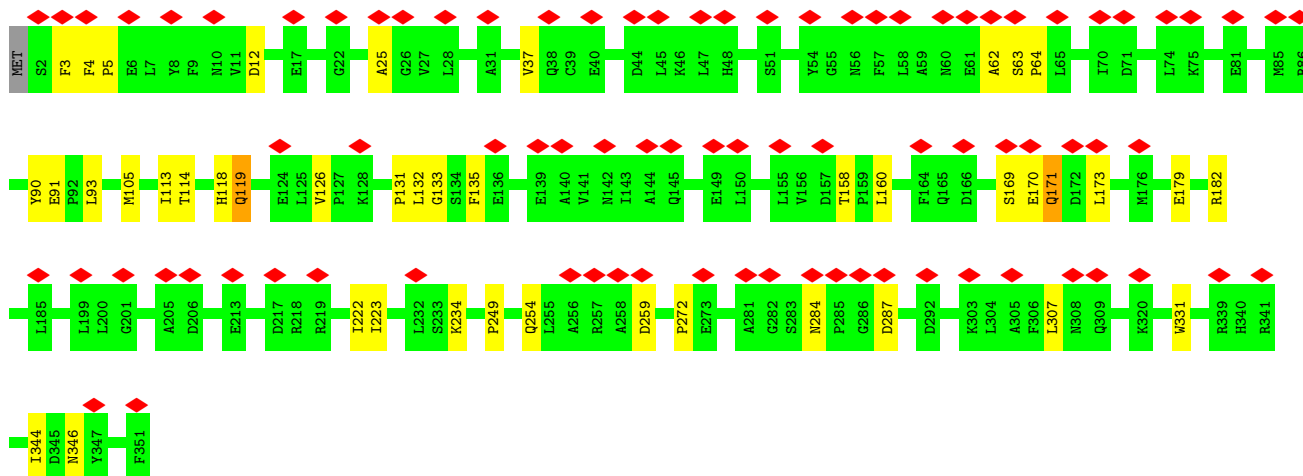


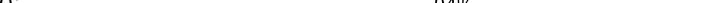
- Molecule 3: V-type proton ATPase 116 kDa subunit a isoform 1

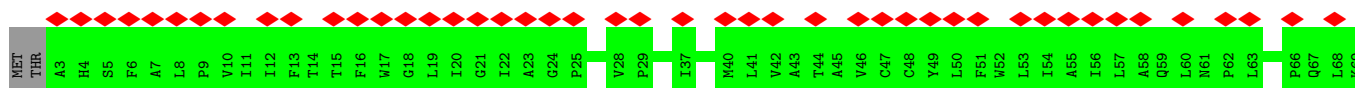




Chain d:

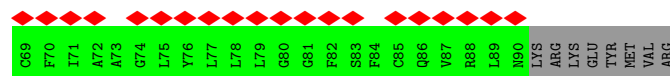
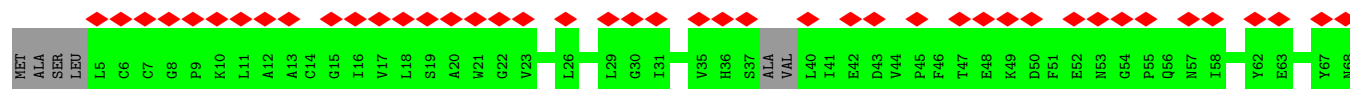
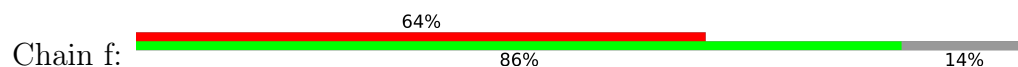


Chain e:  65% 94%

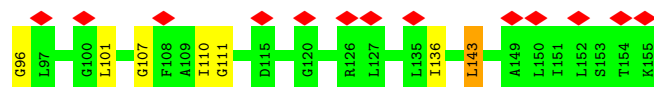
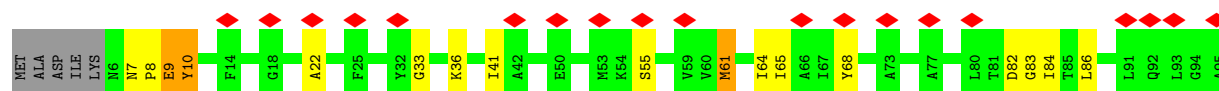
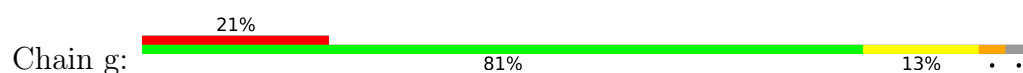




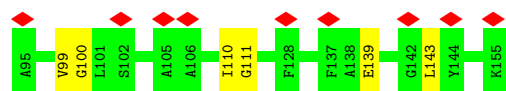
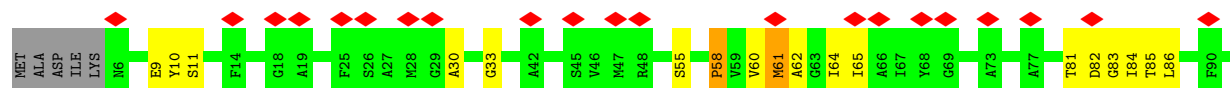
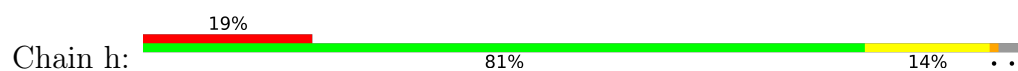
• Molecule 8: Ribonuclease K



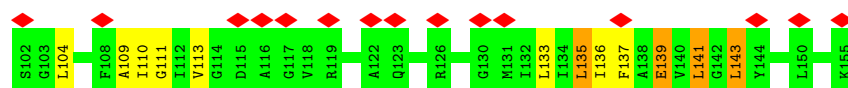
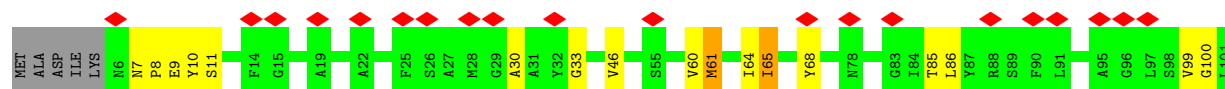
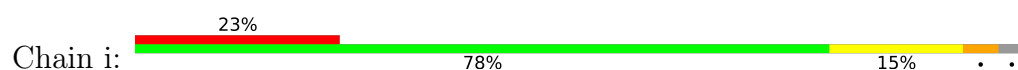
• Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



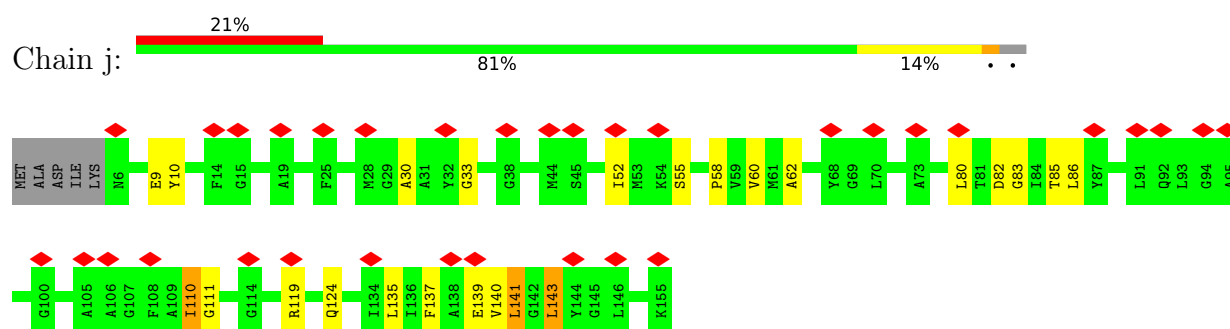
• Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



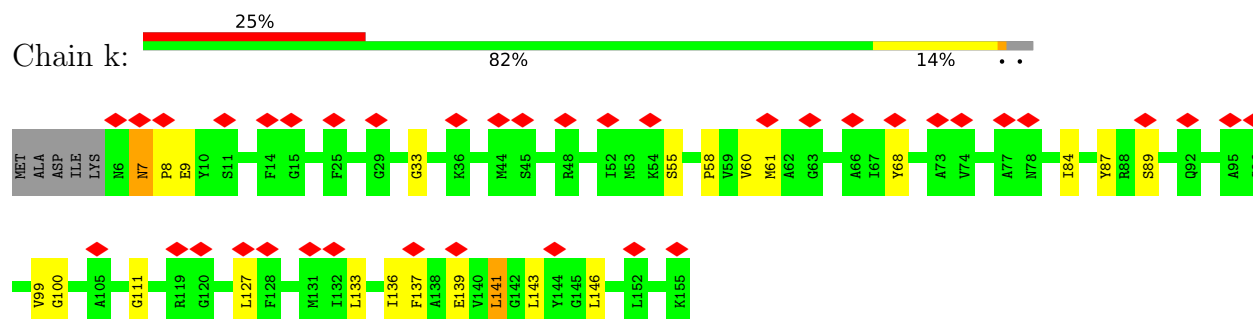
• Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



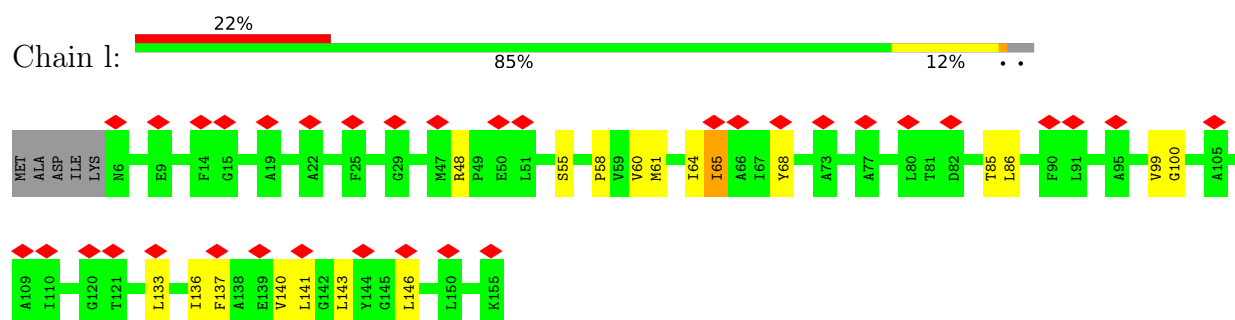
• Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



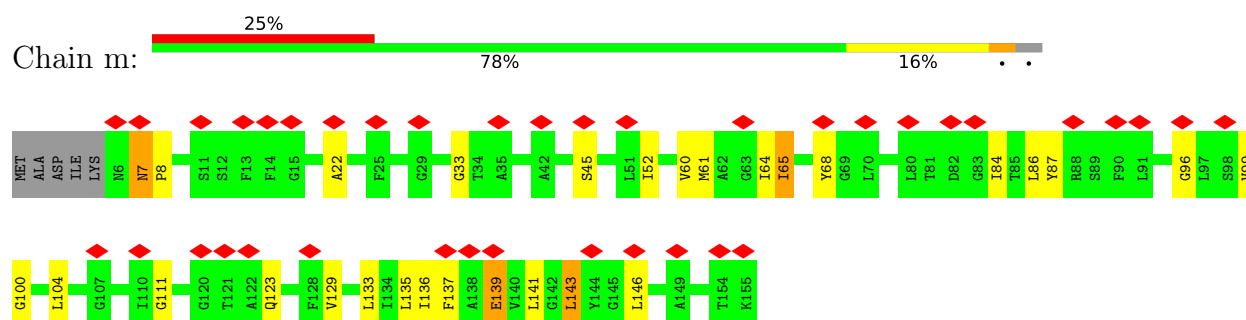
- Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



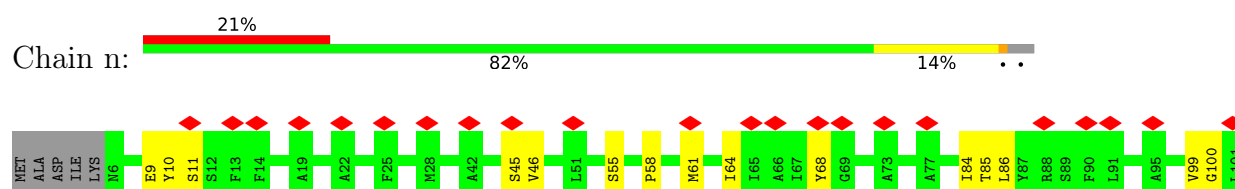
- Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



- Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit



- Molecule 9: V-type proton ATPase 16 kDa proteolipid subunit





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	79654	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 ($e^-/\text{\AA}^2$)	43	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	2.081	Depositor
Minimum map value	-0.649	Depositor
Average map value	-0.002	Depositor
Map value standard deviation	0.056	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	360.4, 360.4, 360.4	wwPDB
Map dimensions	340, 340, 340	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	Depositor

5 Model quality

5.1 Standard geometry

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	H	0.82	1/812 (0.1%)	1.10	4/1091 (0.4%)
2	L	0.85	0/889	1.32	13/1199 (1.1%)
3	a	0.33	0/2170	0.63	1/3014 (0.0%)
4	b	0.45	0/1537	0.68	1/2088 (0.0%)
5	c	0.56	1/347 (0.3%)	0.90	3/466 (0.6%)
6	d	0.59	0/2899	0.70	3/3927 (0.1%)
7	e	0.28	0/383	0.60	0/531
8	f	0.43	0/410	0.81	0/566
9	g	0.53	0/1083	0.82	5/1466 (0.3%)
9	h	0.50	0/1083	0.75	1/1466 (0.1%)
9	i	0.52	0/1083	0.77	4/1466 (0.3%)
9	j	0.47	0/1083	0.68	0/1466
9	k	0.51	0/1083	0.79	3/1466 (0.2%)
9	l	0.49	0/1083	0.75	2/1466 (0.1%)
9	m	0.56	0/1083	0.93	3/1466 (0.2%)
9	n	0.46	0/1083	0.69	0/1466
9	o	0.51	0/1083	0.73	0/1466
10	p	0.48	0/416	0.68	1/571 (0.2%)
All	All	0.53	2/19610 (0.0%)	0.79	44/26647 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	107	PRO	C-O	-5.53	1.17	1.23
5	c	417	PRO	N-CD	5.02	1.54	1.47

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	91	ILE	CA-C-N	7.98	127.74	119.76
2	L	91	ILE	C-N-CA	7.98	127.74	119.76
2	L	108	ARG	CA-C-N	-7.67	112.32	123.05
2	L	108	ARG	C-N-CA	-7.67	112.32	123.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	104	SER	N-CA-C	-7.55	103.06	111.82
9	m	7	ASN	CA-C-N	7.33	127.31	119.76
9	m	7	ASN	C-N-CA	7.33	127.31	119.76
6	d	171	GLN	N-CA-C	-7.25	102.55	111.33
5	c	416	SER	CA-C-N	7.13	126.66	119.24
5	c	416	SER	C-N-CA	7.13	126.66	119.24
2	L	55	PHE	N-CA-C	-6.87	103.88	111.36
5	c	417	PRO	N-CA-C	-6.70	102.92	113.78
6	d	3	PHE	N-CA-C	-6.68	104.70	112.92
9	l	48	ARG	CA-C-N	6.48	126.71	119.32
9	l	48	ARG	C-N-CA	6.48	126.71	119.32
2	L	91	ILE	O-C-N	6.34	128.32	121.10
9	k	7	ASN	CA-C-N	6.29	126.23	119.76
9	k	7	ASN	C-N-CA	6.29	126.23	119.76
2	L	97	PRO	N-CA-C	6.28	120.83	111.41
9	g	10	TYR	N-CA-C	-6.14	104.17	113.51
9	k	9	GLU	CA-C-O	-5.83	115.37	121.55
9	m	129	VAL	N-CA-C	-5.82	104.84	110.72
9	i	9	GLU	CA-C-O	-5.62	115.35	121.87
4	b	94	ILE	N-CA-C	-5.60	105.06	110.72
9	i	7	ASN	CA-C-N	5.57	125.85	119.83
9	i	7	ASN	C-N-CA	5.57	125.85	119.83
9	g	9	GLU	CA-C-O	-5.48	115.51	121.87
3	a	529	LYS	N-CA-C	-5.45	105.45	111.71
10	p	306	VAL	N-CA-C	-5.37	105.30	110.72
9	h	58	PRO	N-CA-C	-5.35	105.11	113.78
2	L	33	HIS	CA-C-N	5.33	126.51	119.84
2	L	33	HIS	C-N-CA	5.33	126.51	119.84
9	g	8	PRO	N-CA-C	-5.33	102.58	111.26
1	H	106	LEU	CA-C-N	5.25	125.41	119.90
1	H	106	LEU	C-N-CA	5.25	125.41	119.90
2	L	105	ILE	CA-C-N	-5.16	112.86	120.28
2	L	105	ILE	C-N-CA	-5.16	112.86	120.28
2	L	105	ILE	N-CA-C	-5.13	105.40	111.00
9	g	7	ASN	CA-C-N	5.06	125.35	119.93
9	g	7	ASN	C-N-CA	5.06	125.35	119.93
1	H	113	HIS	CA-C-N	5.06	127.31	120.38
1	H	113	HIS	C-N-CA	5.06	127.31	120.38
6	d	126	VAL	CB-CA-C	-5.05	109.00	114.35
9	i	10	TYR	N-CA-C	-5.02	105.39	113.28

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	H	803	0	828	5	0
2	L	875	0	883	12	0
3	a	2173	0	962	2	0
4	b	1503	0	1551	15	0
5	c	337	0	338	1	0
6	d	2833	0	2770	32	0
7	e	384	0	176	1	0
8	f	412	0	186	0	0
9	g	1068	0	1136	13	0
9	h	1068	0	1136	18	0
9	i	1068	0	1136	27	0
9	j	1068	0	1136	20	0
9	k	1068	0	1136	19	0
9	l	1068	0	1136	9	0
9	m	1068	0	1136	23	0
9	n	1068	0	1136	15	0
9	o	1068	0	1136	28	0
10	p	406	0	384	3	0
All	All	19338	0	18302	208	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 (208) 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)
9:o:133:LEU:HA	9:o:136:ILE:HD12	1.55	0.89
9:g:61:MET:HA	9:g:64:ILE:HD12	1.59	0.82
9:k:133:LEU:HA	9:k:136:ILE:HD12	1.69	0.74
9:n:133:LEU:HA	9:n:136:ILE:HD12	1.67	0.74
9:n:61:MET:HA	9:n:64:ILE:HD12	1.68	0.74
4:b:59:SER:HB2	9:g:101:LEU:HD23	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:i:61:MET:HE2	9:i:139:GLU:HG2	1.74	0.70
9:j:30:ALA:HB3	9:k:141:LEU:HD21	1.73	0.69
9:h:30:ALA:HB3	9:i:141:LEU:HD21	1.75	0.68
9:i:143:LEU:HD12	9:i:143:LEU:O	1.93	0.67
9:j:143:LEU:HD12	9:j:143:LEU:O	1.95	0.67
9:j:60:VAL:HG22	9:k:137:PHE:CZ	2.33	0.64
9:m:68:TYR:HD1	9:m:146:LEU:HD22	1.63	0.64
9:o:61:MET:N	9:o:61:MET:HE2	2.12	0.64
9:l:61:MET:HA	9:l:64:ILE:HD12	1.78	0.63
9:m:143:LEU:O	9:m:143:LEU:HD12	1.99	0.62
9:o:61:MET:HA	9:o:64:ILE:HD12	1.81	0.62
2:L:62:GLY:O	2:L:86:PRO:HA	1.99	0.61
4:b:168:GLN:HG2	6:d:25:ALA:HB2	1.80	0.61
4:b:132:HIS:CE1	9:o:81:THR:O	2.54	0.61
6:d:12:ASP:OD1	9:g:36:LYS:NZ	2.34	0.61
9:h:62:ALA:HA	9:h:65:ILE:HD12	1.83	0.60
9:i:8:PRO:HB2	9:i:11:SER:OG	2.01	0.60
9:o:68:TYR:HD1	9:o:146:LEU:HD22	1.67	0.59
2:L:85:ILE:N	2:L:86:PRO:HD2	2.17	0.59
9:n:143:LEU:HD12	9:n:147:ILE:HG13	1.85	0.59
2:L:85:ILE:N	2:L:86:PRO:CD	2.66	0.58
6:d:249:PRO:O	9:m:123:GLN:NE2	2.36	0.58
9:g:33:GLY:HA2	9:g:111:GLY:HA3	1.85	0.58
9:h:143:LEU:O	9:h:143:LEU:HD12	2.03	0.57
9:m:61:MET:HE2	9:m:139:GLU:HG2	1.85	0.57
9:g:143:LEU:HD12	9:g:143:LEU:O	2.03	0.57
9:l:60:VAL:HG22	9:m:137:PHE:CZ	2.40	0.57
1:H:107:PRO:HG2	1:H:153:THR:HG21	1.85	0.57
9:k:84:ILE:HG12	9:k:89:SER:HB2	1.85	0.56
9:m:68:TYR:CD1	9:m:146:LEU:HD22	2.39	0.56
6:d:4:PHE:N	6:d:5:PRO:CD	2.67	0.56
6:d:37:VAL:HG13	6:d:344:ILE:HG12	1.88	0.56
9:n:143:LEU:HD12	9:n:143:LEU:O	2.06	0.56
4:b:58:ILE:CD1	4:b:105:ILE:HG12	2.36	0.56
9:i:65:ILE:HA	9:i:68:TYR:CD2	2.42	0.55
9:o:68:TYR:CD1	9:o:146:LEU:HD22	2.42	0.55
4:b:58:ILE:HD11	4:b:105:ILE:HG12	1.89	0.55
6:d:284:ASN:HB2	6:d:287:ASP:HB2	1.89	0.55
9:i:60:VAL:HG22	9:j:137:PHE:CZ	2.41	0.54
6:d:169:SER:C	6:d:171:GLN:H	2.15	0.54
9:l:133:LEU:HA	9:l:136:ILE:HD12	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:k:61:MET:N	9:k:61:MET:HE2	2.23	0.54
9:i:61:MET:HA	9:i:64:ILE:HD12	1.90	0.53
1:H:55:LYS:NZ	2:L:99:ASP:OD2	2.34	0.53
9:m:45:SER:HB2	9:m:52:ILE:HD11	1.90	0.53
9:m:60:VAL:HG22	9:n:137:PHE:CZ	2.43	0.53
9:o:57:ILE:HG23	9:o:61:MET:HE3	1.91	0.52
9:i:46:VAL:HA	9:j:124:GLN:HB2	1.92	0.52
9:o:33:GLY:HA2	9:o:111:GLY:CA	2.40	0.52
4:b:70:TYR:C	4:b:70:TYR:CD1	2.88	0.52
9:h:60:VAL:HG22	9:i:137:PHE:HZ	1.74	0.52
9:n:84:ILE:HG22	9:n:86:LEU:H	1.75	0.52
9:o:9:GLU:O	9:o:11:SER:N	2.44	0.51
6:d:131:PRO:O	6:d:132:LEU:HB2	2.11	0.51
9:j:33:GLY:HA2	9:j:111:GLY:HA3	1.93	0.51
6:d:307:LEU:HA	9:o:43:ALA:HB1	1.92	0.51
9:n:61:MET:HE2	9:n:61:MET:N	2.25	0.51
2:L:96:HIS:CG	2:L:97:PRO:HD2	2.46	0.51
9:g:41:ILE:HG12	9:g:55:SER:HB2	1.93	0.50
9:g:33:GLY:HA2	9:g:111:GLY:CA	2.40	0.50
9:m:65:ILE:HA	9:m:68:TYR:CD2	2.46	0.50
6:d:169:SER:O	6:d:171:GLN:N	2.43	0.50
9:h:65:ILE:HD11	9:h:110:ILE:HG13	1.94	0.50
9:m:65:ILE:HA	9:m:68:TYR:HD2	1.75	0.50
9:j:33:GLY:HA2	9:j:111:GLY:CA	2.42	0.50
6:d:4:PHE:N	6:d:5:PRO:HD2	2.26	0.49
4:b:72:THR:HG23	4:b:156:GLY:HA2	1.94	0.49
9:m:61:MET:HA	9:m:64:ILE:HD12	1.94	0.49
6:d:62:ALA:HB1	6:d:64:PRO:HD2	1.94	0.49
9:n:46:VAL:HG11	9:o:123:GLN:OE1	2.12	0.49
9:n:85:THR:O	9:n:86:LEU:HB3	2.13	0.49
2:L:112:MET:O	2:L:113:PHE:C	2.55	0.49
9:i:60:VAL:HG22	9:j:137:PHE:HZ	1.78	0.49
6:d:169:SER:O	6:d:170:GLU:HB2	2.13	0.48
9:o:86:LEU:O	9:o:89:SER:N	2.47	0.48
9:o:9:GLU:C	9:o:11:SER:H	2.22	0.48
4:b:34:ARG:HH21	5:c:413:GLY:HA2	1.78	0.48
9:o:45:SER:HB2	9:o:52:ILE:HD11	1.96	0.48
9:k:33:GLY:HA2	9:k:111:GLY:HA3	1.96	0.48
9:n:9:GLU:O	9:n:11:SER:N	2.46	0.48
9:o:86:LEU:O	9:o:87:TYR:C	2.55	0.48
2:L:62:GLY:HA2	2:L:85:ILE:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:h:60:VAL:HG22	9:i:137:PHE:CZ	2.49	0.47
9:o:65:ILE:HA	9:o:68:TYR:HD2	1.78	0.47
4:b:196:LEU:HD22	9:o:78:ASN:OD1	2.15	0.47
9:i:61:MET:HA	9:i:64:ILE:CD1	2.46	0.46
4:b:196:LEU:CD1	9:o:74:VAL:HG13	2.46	0.46
6:d:105:MET:HG2	6:d:133:GLY:HA2	1.96	0.46
9:h:60:VAL:HA	9:i:137:PHE:CE2	2.50	0.46
9:i:8:PRO:HG2	9:i:86:LEU:HD22	1.96	0.46
9:m:133:LEU:HA	9:m:136:ILE:HD12	1.97	0.46
6:d:158:THR:HG22	6:d:160:LEU:H	1.80	0.46
6:d:331:TRP:CE2	6:d:346:ASN:HB2	2.51	0.46
9:j:60:VAL:HG22	9:k:137:PHE:HZ	1.77	0.46
9:i:33:GLY:HA2	9:i:111:GLY:HA3	1.97	0.46
6:d:63:SER:N	6:d:64:PRO:HD2	2.31	0.46
9:m:133:LEU:O	9:m:136:ILE:HB	2.16	0.46
9:l:68:TYR:CD1	9:l:146:LEU:HD22	2.51	0.46
4:b:168:GLN:HG2	6:d:25:ALA:CB	2.45	0.45
9:g:107:GLY:HA2	9:g:110:ILE:HD12	1.97	0.45
9:i:46:VAL:HG21	10:p:338:ILE:HG21	1.98	0.45
9:k:33:GLY:HA2	9:k:111:GLY:CA	2.46	0.45
9:m:84:ILE:HG22	9:m:86:LEU:H	1.81	0.45
9:g:84:ILE:HG22	9:g:86:LEU:H	1.82	0.45
3:a:403:PHE:O	3:a:407:PHE:N	2.50	0.45
9:j:52:ILE:CD1	9:k:127:LEU:HA	2.46	0.45
9:i:30:ALA:HB3	9:j:141:LEU:HD21	1.99	0.45
9:j:58:PRO:HG3	9:j:135:LEU:HD11	1.99	0.45
6:d:118:HIS:O	6:d:119:GLN:C	2.60	0.45
6:d:331:TRP:CZ2	6:d:346:ASN:HB2	2.52	0.45
9:m:65:ILE:O	9:m:68:TYR:HB2	2.16	0.45
9:o:133:LEU:O	9:o:136:ILE:HB	2.17	0.45
9:n:68:TYR:CD1	9:n:146:LEU:HD22	2.52	0.44
4:b:190:GLY:HA2	9:o:23:MET:HG2	1.99	0.44
9:h:55:SER:O	9:h:58:PRO:HD2	2.18	0.44
9:m:33:GLY:HA2	9:m:111:GLY:HA3	1.99	0.44
6:d:234:LYS:HE2	6:d:259:ASP:OD1	2.18	0.44
9:j:62:ALA:HA	9:j:110:ILE:HD12	2.00	0.44
9:k:61:MET:HE2	9:k:61:MET:CA	2.47	0.44
9:m:99:VAL:HG13	9:m:100:GLY:N	2.33	0.44
9:h:9:GLU:O	9:h:11:SER:N	2.51	0.44
9:k:99:VAL:HG13	9:k:100:GLY:N	2.33	0.44
4:b:59:SER:O	4:b:63:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:h:9:GLU:HG3	9:h:10:TYR:N	2.33	0.44
2:L:34:PRO:HB2	2:L:36:PHE:CE2	2.53	0.44
9:m:22:ALA:HB2	9:m:96:GLY:HA2	1.99	0.44
9:o:65:ILE:N	9:o:65:ILE:CD1	2.81	0.44
9:o:107:GLY:HA2	9:o:110:ILE:HD12	1.99	0.44
9:j:80:LEU:C	9:k:87:TYR:OH	2.61	0.43
9:o:65:ILE:HA	9:o:68:TYR:CD2	2.52	0.43
9:o:84:ILE:HG22	9:o:86:LEU:H	1.82	0.43
9:m:104:LEU:C	9:m:104:LEU:HD23	2.43	0.43
6:d:169:SER:C	6:d:171:GLN:N	2.76	0.43
9:g:65:ILE:O	9:g:68:TYR:HB2	2.18	0.43
6:d:113:ILE:O	6:d:114:THR:HB	2.19	0.43
9:i:104:LEU:HD23	9:i:104:LEU:C	2.43	0.43
9:i:143:LEU:HD12	9:i:143:LEU:C	2.39	0.43
2:L:61:ILE:HG22	2:L:62:GLY:N	2.33	0.43
9:i:61:MET:SD	9:i:135:LEU:HB3	2.59	0.43
4:b:58:ILE:HG23	4:b:59:SER:N	2.34	0.43
9:g:82:ASP:OD1	9:g:83:GLY:N	2.51	0.43
9:h:61:MET:HA	9:h:64:ILE:CD1	2.49	0.43
2:L:6:LYS:NZ	2:L:59:ASP:O	2.48	0.43
9:m:61:MET:SD	9:m:135:LEU:HB3	2.59	0.43
9:k:60:VAL:HG22	9:l:137:PHE:CZ	2.54	0.42
9:n:99:VAL:HG13	9:n:100:GLY:N	2.34	0.42
9:j:80:LEU:C	9:k:87:TYR:HH	2.27	0.42
9:j:85:THR:O	9:j:86:LEU:HB3	2.19	0.42
4:b:67:TRP:HA	4:b:70:TYR:CE2	2.54	0.42
9:h:61:MET:HA	9:h:64:ILE:HD12	2.00	0.42
9:i:65:ILE:O	9:i:68:TYR:HB2	2.19	0.42
9:o:48:ARG:HB2	9:o:50:GLU:CD	2.45	0.42
3:a:407:PHE:O	3:a:409:ASP:N	2.53	0.42
9:j:9:GLU:C	9:j:10:TYR:CD2	2.98	0.42
10:p:336:ASP:C	10:p:336:ASP:OD1	2.62	0.42
9:j:119:ARG:HH11	10:p:339:ILE:HD12	1.85	0.42
9:k:55:SER:O	9:k:58:PRO:HD2	2.19	0.42
9:i:8:PRO:O	9:i:11:SER:OG	2.36	0.42
9:j:82:ASP:OD1	9:j:83:GLY:N	2.53	0.42
9:l:99:VAL:HG13	9:l:100:GLY:N	2.34	0.42
9:h:30:ALA:CB	9:i:141:LEU:HD21	2.46	0.42
9:o:55:SER:O	9:o:58:PRO:HD2	2.20	0.42
6:d:133:GLY:O	6:d:135:PHE:N	2.53	0.41
9:h:33:GLY:HA2	9:h:111:GLY:CA	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:n:55:SER:O	9:n:58:PRO:HD2	2.20	0.41
9:g:9:GLU:O	9:g:10:TYR:HB2	2.19	0.41
9:h:85:THR:O	9:h:86:LEU:HB3	2.19	0.41
9:i:99:VAL:HG13	9:i:100:GLY:N	2.35	0.41
9:j:55:SER:O	9:j:58:PRO:HD2	2.20	0.41
9:k:7:ASN:HB3	9:k:8:PRO:HD2	2.01	0.41
9:m:7:ASN:HA	9:m:8:PRO:HD3	1.94	0.41
9:m:86:LEU:O	9:m:87:TYR:C	2.63	0.41
9:n:45:SER:HB2	9:o:127:LEU:HD21	2.03	0.41
2:L:13:ASP:OD1	2:L:13:ASP:C	2.64	0.41
9:l:85:THR:O	9:l:86:LEU:HB3	2.20	0.41
6:d:254:GLN:OE1	6:d:254:GLN:N	2.49	0.41
7:e:76:VAL:O	7:e:77:ARG:C	2.63	0.41
9:i:133:LEU:HA	9:i:136:ILE:HD12	2.02	0.41
9:o:99:VAL:HG13	9:o:100:GLY:N	2.35	0.41
1:H:88:ASN:OD1	1:H:89:LYS:N	2.52	0.41
6:d:170:GLU:O	6:d:173:LEU:N	2.53	0.41
6:d:179:GLU:O	6:d:182:ARG:HG2	2.21	0.41
6:d:272:PRO:HD2	9:n:123:GLN:HG2	2.03	0.41
9:h:99:VAL:HG13	9:h:100:GLY:N	2.35	0.41
9:i:109:ALA:O	9:i:113:VAL:HG22	2.21	0.41
9:k:133:LEU:O	9:k:136:ILE:HB	2.21	0.41
6:d:90:TYR:O	6:d:91:GLU:C	2.62	0.41
6:d:93:LEU:C	6:d:93:LEU:HD23	2.45	0.41
6:d:131:PRO:O	6:d:132:LEU:CB	2.69	0.41
9:k:68:TYR:CD1	9:k:146:LEU:HD22	2.56	0.41
9:l:55:SER:O	9:l:58:PRO:HD2	2.21	0.41
9:m:61:MET:HA	9:m:64:ILE:CD1	2.51	0.41
1:H:69:LEU:C	1:H:69:LEU:HD23	2.46	0.40
2:L:85:ILE:HB	2:L:86:PRO:HD3	2.03	0.40
9:h:82:ASP:OD1	9:h:83:GLY:N	2.51	0.40
9:h:81:THR:H	9:h:84:ILE:HD11	1.87	0.40
9:k:61:MET:HE2	9:k:61:MET:HA	2.03	0.40
1:H:122:THR:O	1:H:126:ARG:HA	2.22	0.40
6:d:222:ILE:HG23	6:d:223:ILE:N	2.36	0.40
9:i:85:THR:O	9:i:86:LEU:HB3	2.21	0.40
9:l:65:ILE:HA	9:l:68:TYR:CD2	2.56	0.40
9:g:22:ALA:HB2	9:g:96:GLY:HA2	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	H	102/247 (41%)	97 (95%)	5 (5%)	0	100	100
2	L	108/119 (91%)	101 (94%)	5 (5%)	2 (2%)	6	32
3	a	435/838 (52%)	426 (98%)	9 (2%)	0	100	100
4	b	201/205 (98%)	198 (98%)	3 (2%)	0	100	100
5	c	39/463 (8%)	39 (100%)	0	0	100	100
6	d	348/351 (99%)	329 (94%)	18 (5%)	1 (0%)	36	71
7	e	76/81 (94%)	75 (99%)	1 (1%)	0	100	100
8	f	80/98 (82%)	77 (96%)	3 (4%)	0	100	100
9	g	148/155 (96%)	146 (99%)	2 (1%)	0	100	100
9	h	148/155 (96%)	145 (98%)	3 (2%)	0	100	100
9	i	148/155 (96%)	145 (98%)	3 (2%)	0	100	100
9	j	148/155 (96%)	146 (99%)	2 (1%)	0	100	100
9	k	148/155 (96%)	145 (98%)	3 (2%)	0	100	100
9	l	148/155 (96%)	145 (98%)	3 (2%)	0	100	100
9	m	148/155 (96%)	145 (98%)	3 (2%)	0	100	100
9	n	148/155 (96%)	145 (98%)	2 (1%)	1 (1%)	18	55
9	o	148/155 (96%)	143 (97%)	4 (3%)	1 (1%)	18	55
10	p	50/350 (14%)	46 (92%)	3 (6%)	1 (2%)	6	31
All	All	2771/4147 (67%)	2693 (97%)	72 (3%)	6 (0%)	44	77

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	p	300	ASN
9	n	10	TYR
9	o	10	TYR
2	L	100	ALA

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Mol	Chain	Res	Type
6	d	119	GLN
2	L	61	ILE

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	H	84/212 (40%)	84 (100%)	0	100	100
2	L	94/100 (94%)	94 (100%)	0	100	100
4	b	156/158 (99%)	156 (100%)	0	100	100
5	c	36/395 (9%)	36 (100%)	0	100	100
6	d	305/306 (100%)	305 (100%)	0	100	100
9	g	109/113 (96%)	106 (97%)	3 (3%)	38	59
9	h	109/113 (96%)	107 (98%)	2 (2%)	51	67
9	i	109/113 (96%)	102 (94%)	7 (6%)	16	38
9	j	109/113 (96%)	104 (95%)	5 (5%)	24	46
9	k	109/113 (96%)	106 (97%)	3 (3%)	38	59
9	l	109/113 (96%)	105 (96%)	4 (4%)	30	51
9	m	109/113 (96%)	105 (96%)	4 (4%)	30	51
9	n	109/113 (96%)	107 (98%)	2 (2%)	51	67
9	o	109/113 (96%)	104 (95%)	5 (5%)	24	46
10	p	41/313 (13%)	41 (100%)	0	100	100
All	All	1697/2501 (68%)	1662 (98%)	35 (2%)	46	65

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	g	61	MET
9	g	136	ILE
9	g	143	LEU
9	h	61	MET

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Mol	Chain	Res	Type
9	h	139	GLU
9	i	61	MET
9	i	65	ILE
9	i	110	ILE
9	i	135	LEU
9	i	139	GLU
9	i	141	LEU
9	i	143	LEU
9	j	110	ILE
9	j	139	GLU
9	j	140	VAL
9	j	141	LEU
9	j	143	LEU
9	k	139	GLU
9	k	141	LEU
9	k	143	LEU
9	l	65	ILE
9	l	140	VAL
9	l	141	LEU
9	l	143	LEU
9	m	65	ILE
9	m	139	GLU
9	m	141	LEU
9	m	143	LEU
9	n	139	GLU
9	n	143	LEU
9	o	65	ILE
9	o	139	GLU
9	o	140	VAL
9	o	141	LEU
9	o	143	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	85	GLN
1	H	137	ASN
6	d	30	GLN
6	d	88	HIS
9	m	124	GLN
10	p	308	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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.

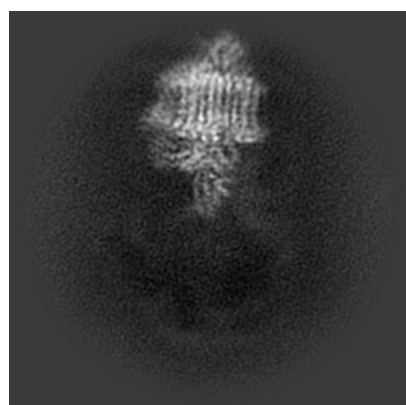
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21350. These allow visual inspection of the internal detail of the map and identification of artifacts.

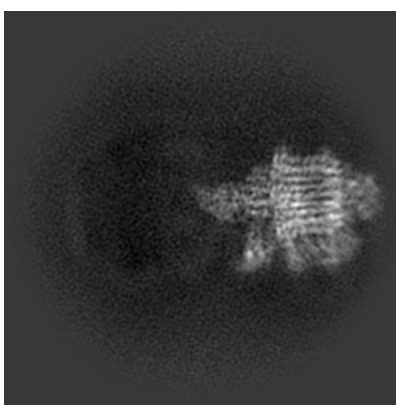
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

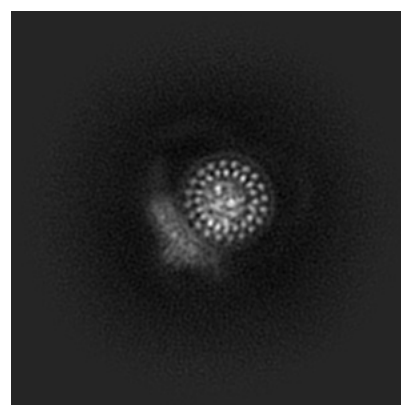
6.1.1 Primary map



X



Y

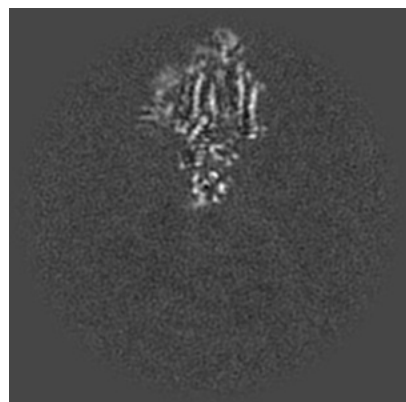


Z

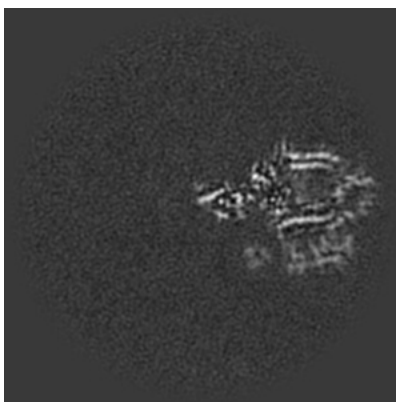
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

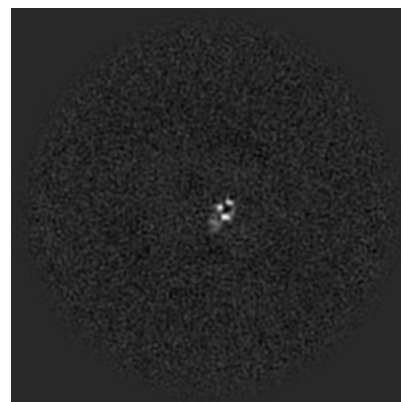
6.2.1 Primary map



X Index: 170



Y Index: 170

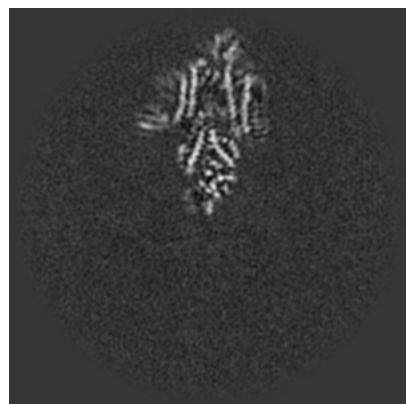


Z Index: 170

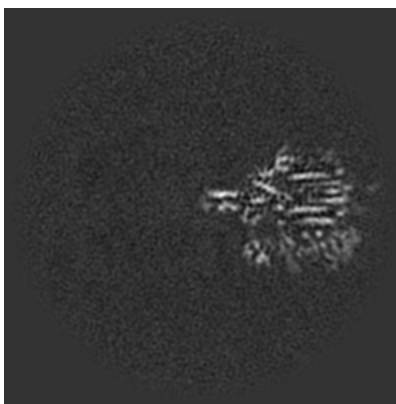
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

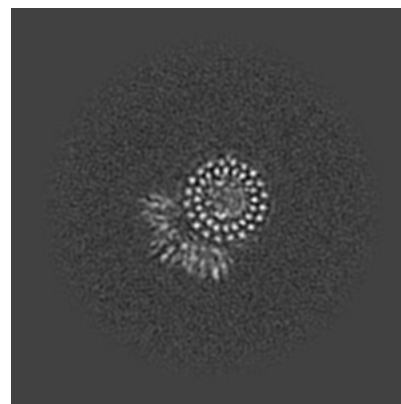
6.3.1 Primary map



X Index: 176



Y Index: 158

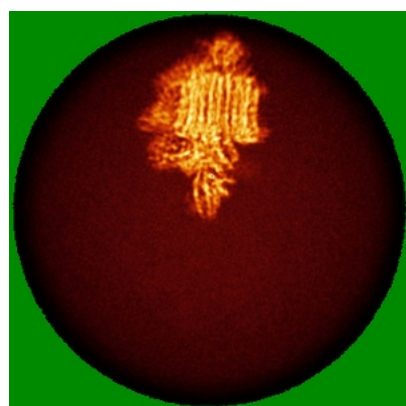


Z Index: 248

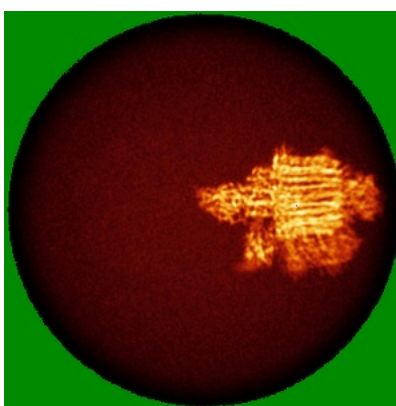
The images above show the largest variance slices of the map in three orthogonal directions.

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

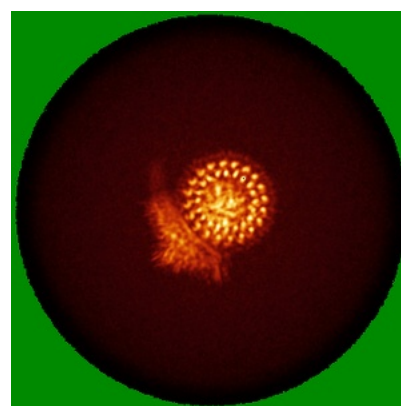
6.4.1 Primary map



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.

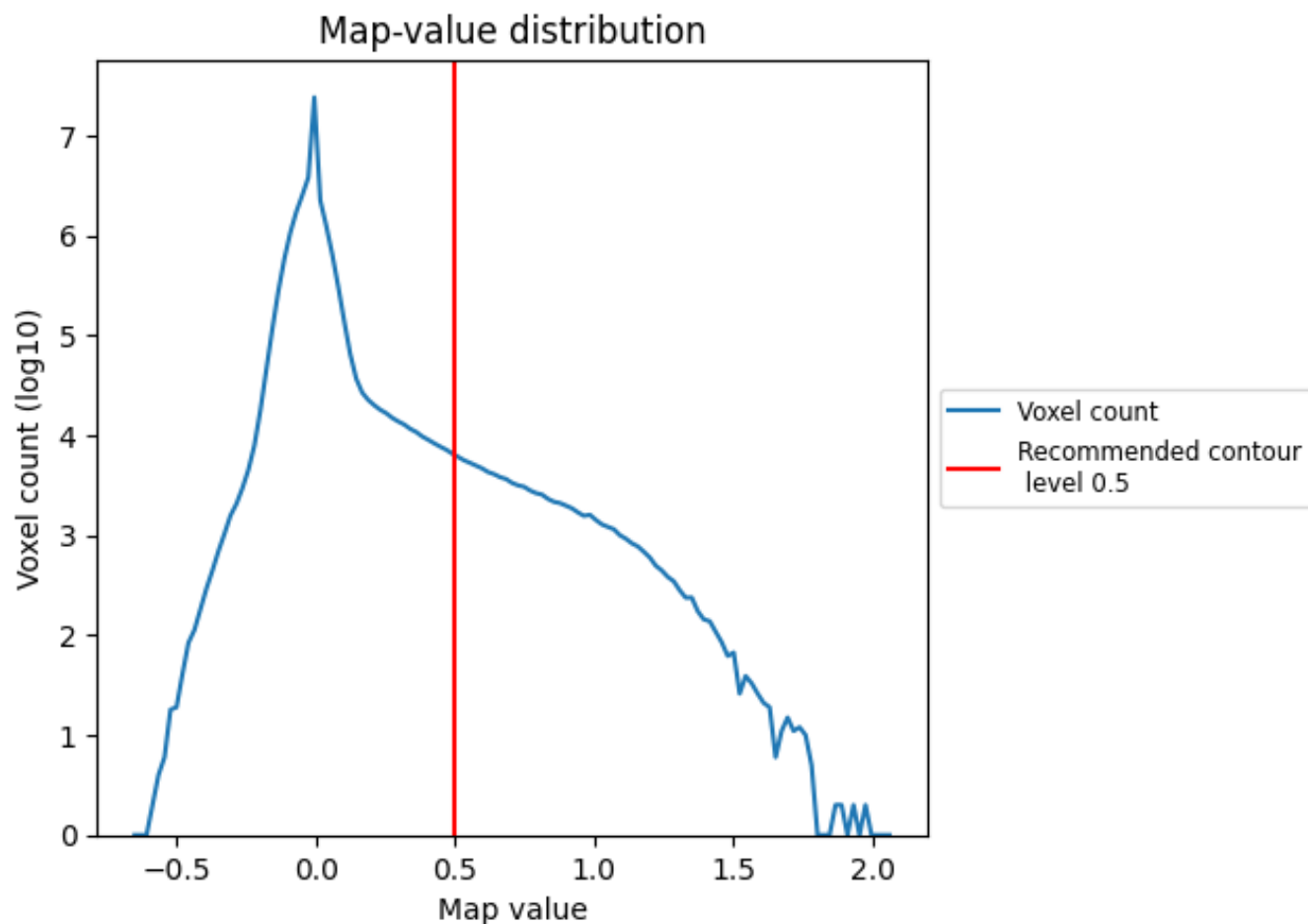
6.6 Mask visualisation

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

7 Map analysis [i](#)

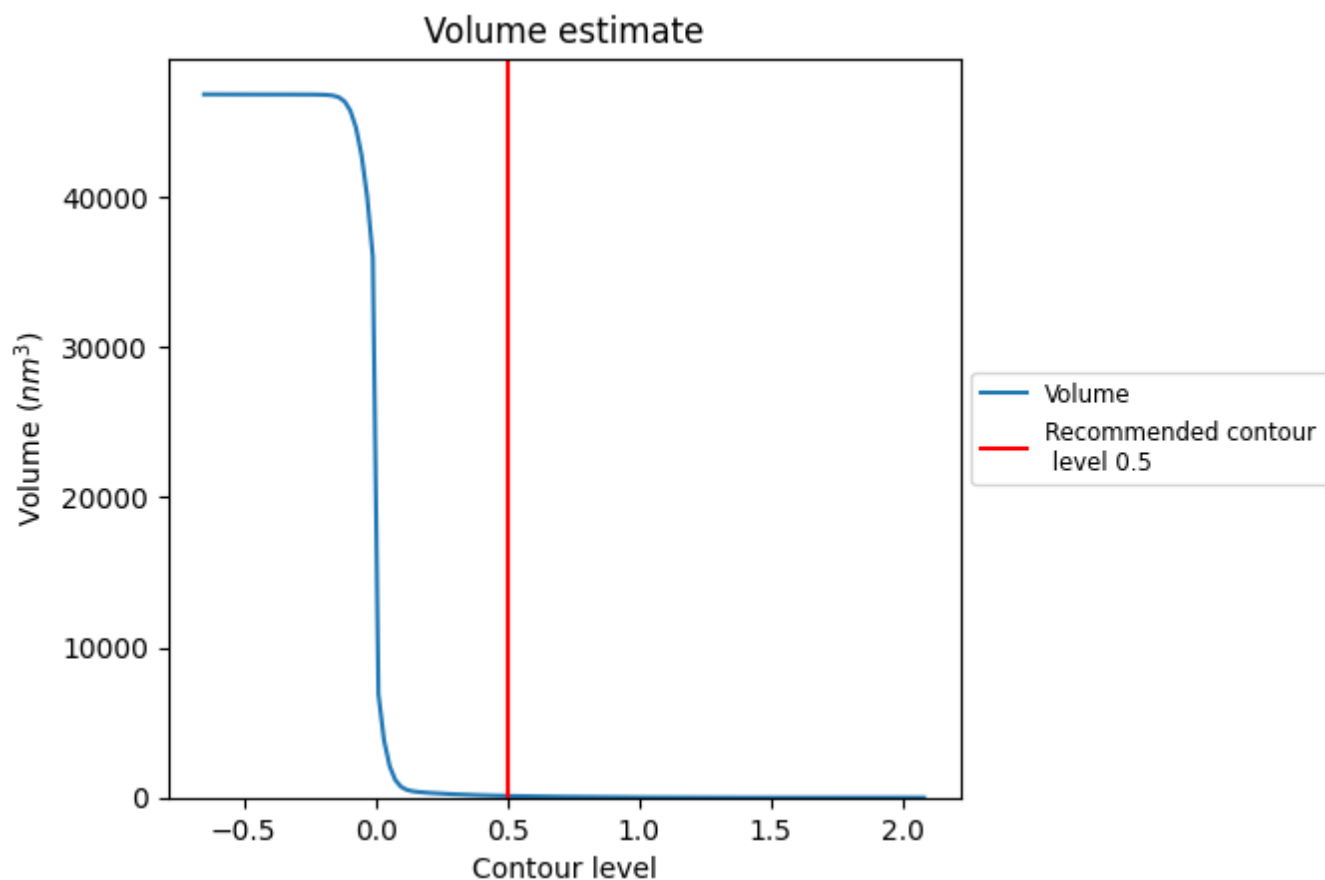
This section contains the results of statistical analysis of the map.

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.

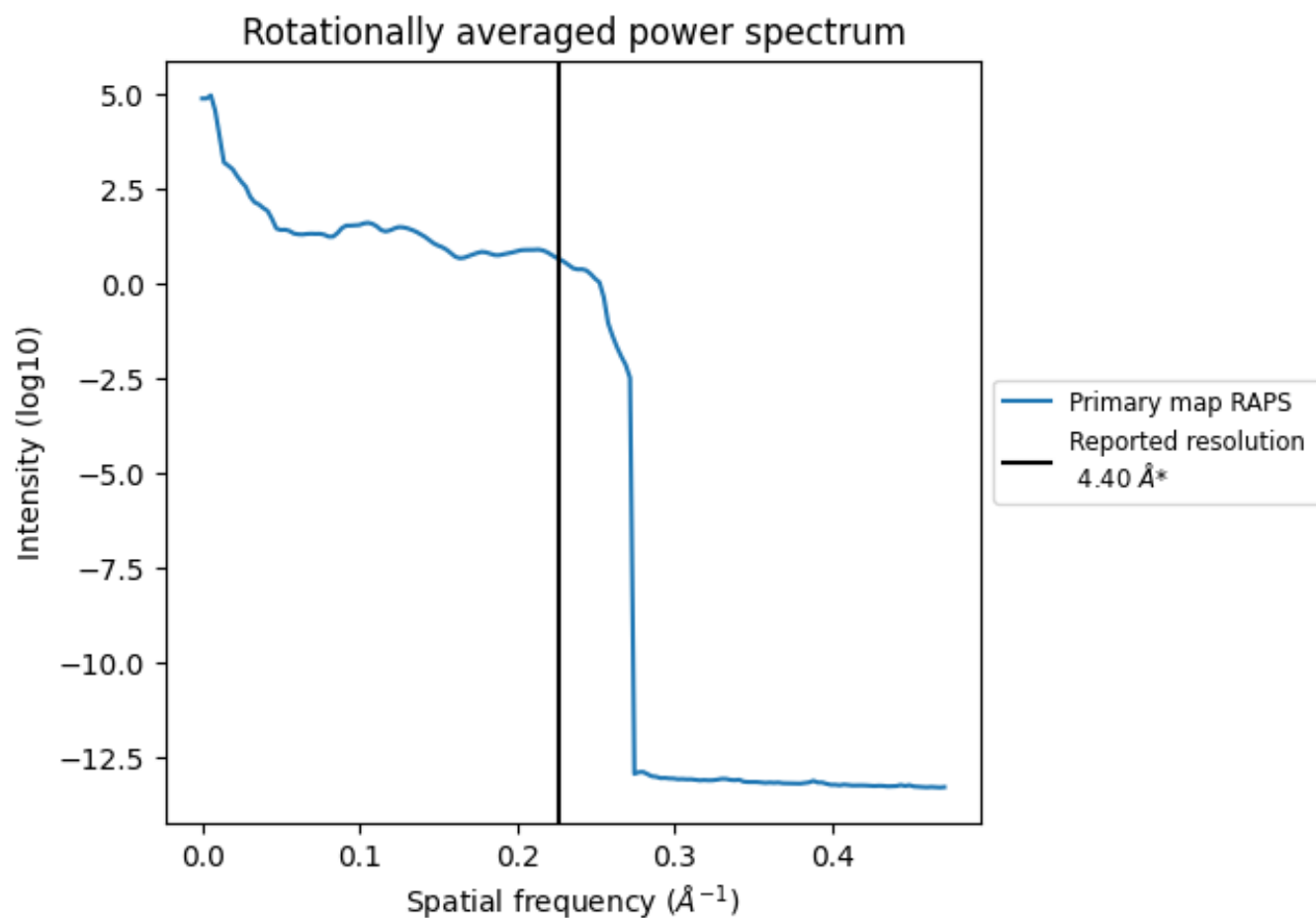
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 110 nm³; this corresponds to an approximate mass of 99 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 0.227 Å⁻¹

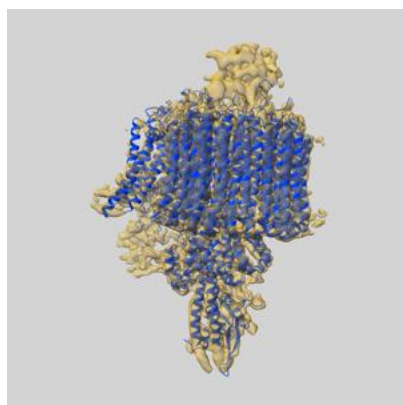
8 Fourier-Shell correlation ⓘ

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

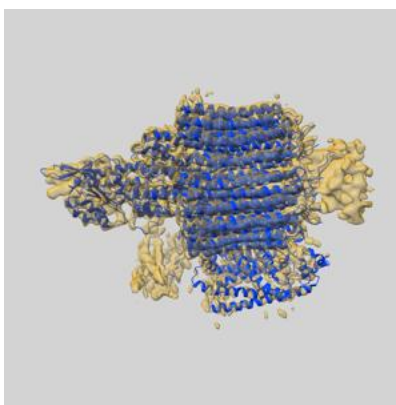
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-21350 and PDB model 6VQH. Per-residue inclusion information can be found in section [3](#) on page [6](#).

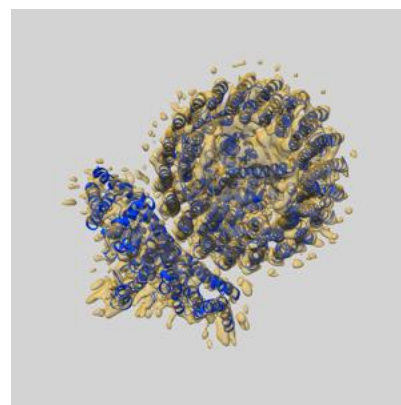
9.1 Map-model overlay [i](#)



X



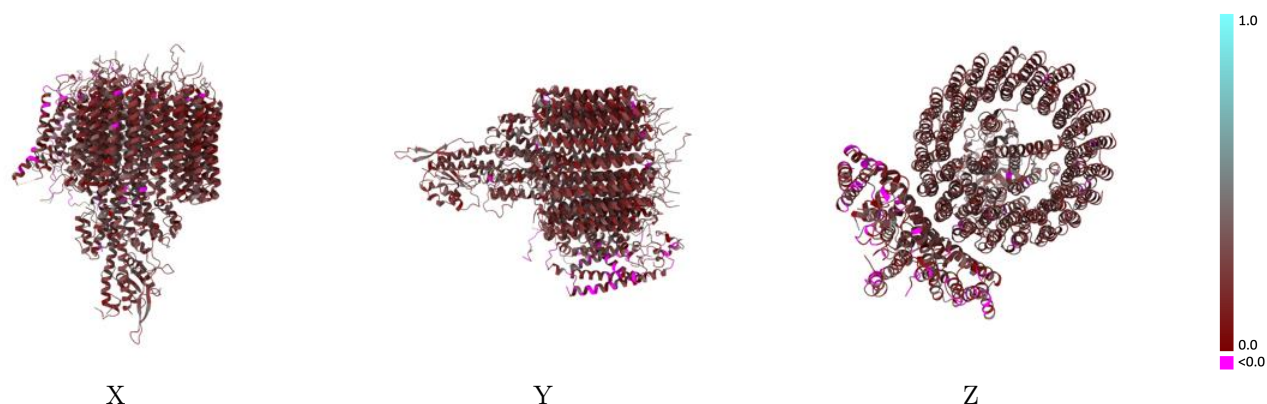
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.5 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](#)

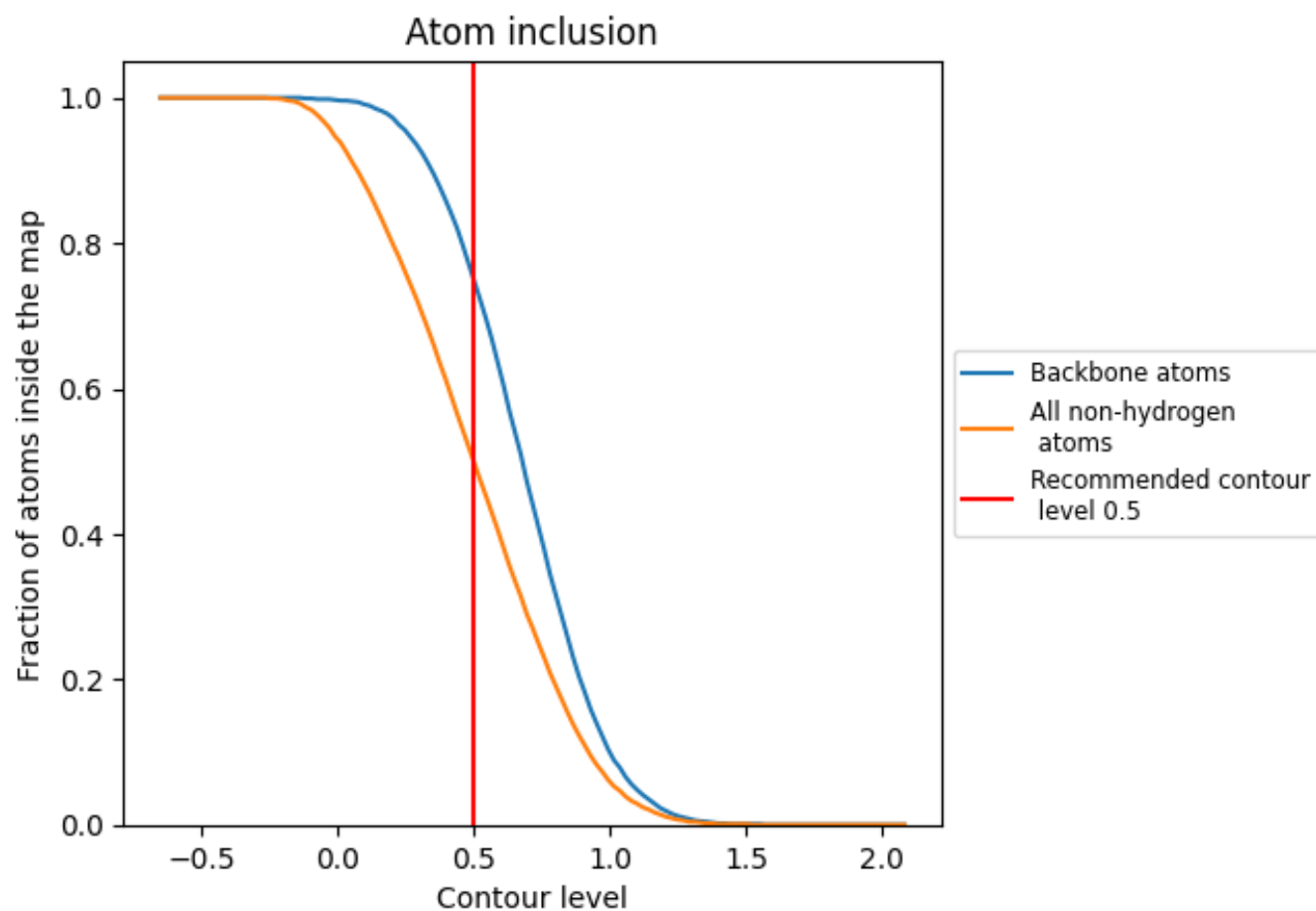


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 50% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.4990	 0.2420
H	 0.5810	 0.2700
L	 0.5440	 0.2710
a	 0.4050	 0.1940
b	 0.5020	 0.2460
c	 0.4530	 0.2970
d	 0.5250	 0.2850
e	 0.3700	 0.1980
f	 0.2860	 0.1670
g	 0.5250	 0.2440
h	 0.5160	 0.2430
i	 0.5150	 0.2370
j	 0.5160	 0.2280
k	 0.4980	 0.2230
l	 0.5120	 0.2370
m	 0.5230	 0.2360
n	 0.5370	 0.2420
o	 0.5010	 0.2390
p	 0.5390	 0.2690

