



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 04:47 PM UTC

PDB ID : 6E2F / pdb_00006e2f
EMDB ID : EMD-8961
Title : Cryo-EM structure of human TRPV6 in complex with Calmodulin
Authors : Singh, A.K.; McGoldrick, L.L.; Sobolevsky, A.I.
Deposited on : 2018-07-11
Resolution : 3.90 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:

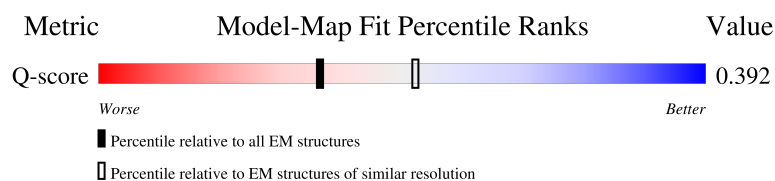
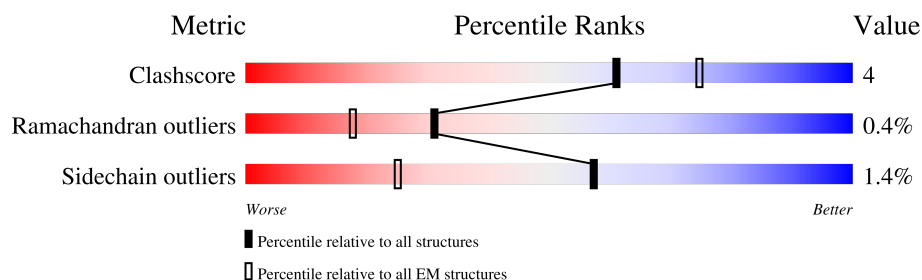
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.90 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	8855 (3.40 - 4.40)

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	725	
1	B	725	
1	C	725	
1	D	725	

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Mol	Chain	Length	Quality of chain
2	E	149	41% 88% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21204 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 Transient receptor potential cation channel subfamily V member 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	613	Total	C	N	O	S	0	0
			4919	3178	833	868	40		
1	B	619	Total	C	N	O	S	0	0
			4978	3212	850	876	40		
1	C	647	Total	C	N	O	S	0	0
			5217	3355	905	917	40		
1	D	613	Total	C	N	O	S	0	0
			4919	3178	833	868	40		

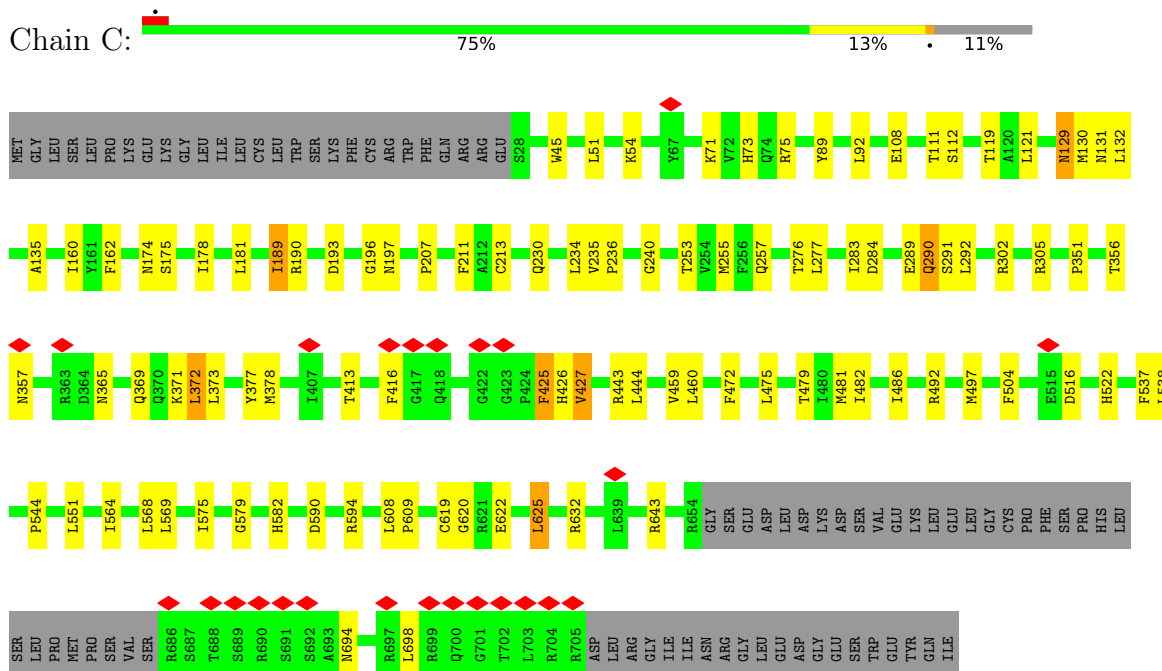
- Molecule 2 is a protein called Calmodulin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	148	Total	C	N	O	S	0	0
			1165	714	188	254	9		

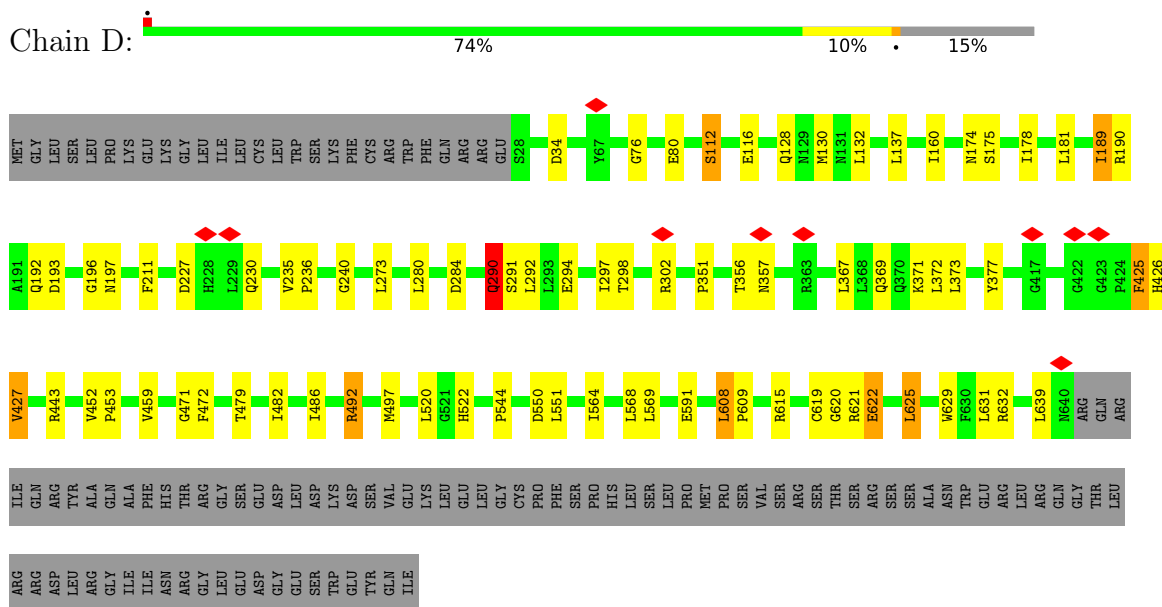
- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

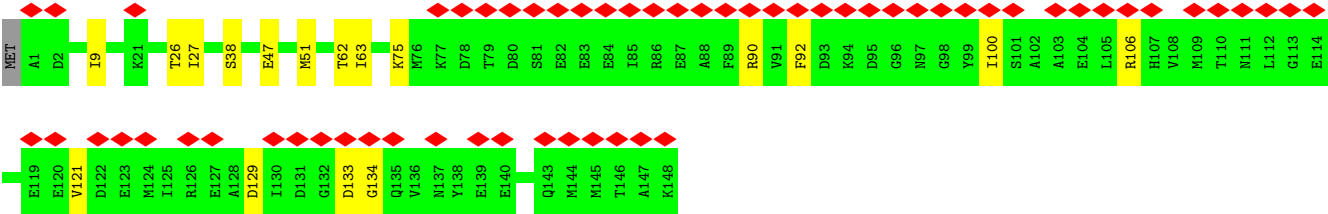
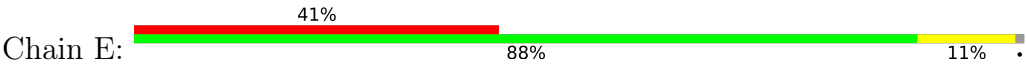
- Molecule 1: Transient receptor potential cation channel subfamily V member 6



- Molecule 1: Transient receptor potential cation channel subfamily V member 6



● Molecule 2: Calmodulin-1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	180600	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 ($e^-/\text{\AA}^2$)	55.0	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	1.381	Depositor
Minimum map value	-0.664	Depositor
Average map value	0.014	Depositor
Map value standard deviation	0.074	Depositor
Recommended contour level	0.34	Depositor
Map size (Å)	270.0, 270.0, 270.0	wwPDB
Map dimensions	250, 250, 250	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	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

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	0.49	1/5032 (0.0%)	0.85	18/6828 (0.3%)
1	B	0.47	0/5091	0.82	11/6905 (0.2%)
1	C	0.47	0/5334	0.82	13/7229 (0.2%)
1	D	0.48	1/5032 (0.0%)	0.84	13/6828 (0.2%)
2	E	0.27	0/1177	0.72	2/1580 (0.1%)
All	All	0.47	2/21666 (0.0%)	0.83	57/29370 (0.2%)

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	3
1	B	0	3
1	C	0	4
1	D	0	3
All	All	0	13

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	550	ASP	C-N	-8.93	1.21	1.33
1	D	550	ASP	C-N	-5.70	1.25	1.33

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	427	VAL	N-CA-C	-11.70	102.60	113.71
1	D	427	VAL	N-CA-C	-11.49	102.80	113.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	427	VAL	N-CA-C	-10.68	103.56	113.71
1	B	550	ASP	CA-C-N	7.63	132.28	120.68
1	B	550	ASP	C-N-CA	7.63	132.28	120.68
1	A	369	GLN	CA-C-N	-6.60	112.00	122.17
1	A	369	GLN	C-N-CA	-6.60	112.00	122.17
1	D	622	GLU	N-CA-C	-6.57	104.30	112.90
1	A	622	GLU	N-CA-C	-6.44	104.46	112.90
1	A	46	GLU	CA-C-N	-6.43	111.80	121.83
1	A	46	GLU	C-N-CA	-6.43	111.80	121.83
1	C	622	GLU	N-CA-C	-6.43	104.48	112.90
1	A	550	ASP	CA-C-N	6.26	130.18	120.60
1	A	550	ASP	C-N-CA	6.26	130.18	120.60
1	C	369	GLN	CA-C-N	-6.23	112.58	122.17
1	C	369	GLN	C-N-CA	-6.23	112.58	122.17
1	B	622	GLU	N-CA-C	-6.13	105.62	113.23
1	D	369	GLN	CA-C-N	-6.09	112.79	122.17
1	D	369	GLN	C-N-CA	-6.09	112.79	122.17
1	C	290	GLN	CA-C-N	5.78	132.58	121.54
1	C	290	GLN	C-N-CA	5.78	132.58	121.54
1	D	472	PHE	CA-C-N	5.73	132.49	121.54
1	D	472	PHE	C-N-CA	5.73	132.49	121.54
1	D	112	SER	CA-C-N	5.73	132.49	121.54
1	D	112	SER	C-N-CA	5.73	132.49	121.54
2	E	75	LYS	CA-C-N	5.72	129.84	121.31
2	E	75	LYS	C-N-CA	5.72	129.84	121.31
1	B	112	SER	CA-C-N	5.72	132.47	121.54
1	B	112	SER	C-N-CA	5.72	132.47	121.54
1	B	369	GLN	CA-C-N	-5.69	113.41	122.17
1	B	369	GLN	C-N-CA	-5.69	113.41	122.17
1	C	472	PHE	CA-C-N	5.67	132.36	121.54
1	C	472	PHE	C-N-CA	5.67	132.36	121.54
1	A	472	PHE	CA-C-N	5.62	132.27	121.54
1	A	472	PHE	C-N-CA	5.62	132.27	121.54
1	A	289	GLU	CA-C-N	5.60	132.24	121.54
1	A	289	GLU	C-N-CA	5.60	132.24	121.54
1	D	290	GLN	CA-C-N	5.56	132.16	121.54
1	D	290	GLN	C-N-CA	5.56	132.16	121.54
1	B	289	GLU	CA-C-N	5.39	131.84	121.54
1	B	289	GLU	C-N-CA	5.39	131.84	121.54
1	C	289	GLU	CA-C-N	5.35	131.76	121.54
1	C	289	GLU	C-N-CA	5.35	131.76	121.54
1	D	639	LEU	CA-CB-CG	5.34	134.98	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	372	LEU	CA-CB-CG	5.24	134.65	116.30
1	A	369	GLN	N-CA-CB	-5.21	102.88	111.01
1	B	471	GLY	CA-C-N	5.12	131.32	121.54
1	B	471	GLY	C-N-CA	5.12	131.32	121.54
1	A	112	SER	CA-C-N	5.11	131.30	121.54
1	A	112	SER	C-N-CA	5.11	131.30	121.54
1	A	471	GLY	CA-C-N	5.11	131.29	121.54
1	A	471	GLY	C-N-CA	5.11	131.29	121.54
1	A	372	LEU	CA-CB-CG	5.09	134.10	116.30
1	C	112	SER	CA-C-N	5.06	131.21	121.54
1	C	112	SER	C-N-CA	5.06	131.21	121.54
1	D	471	GLY	CA-C-N	5.04	131.17	121.54
1	D	471	GLY	C-N-CA	5.04	131.17	121.54

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	290	GLN	Peptide
1	A	371	LYS	Peptide
1	A	425	PHE	Peptide
1	B	290	GLN	Peptide
1	B	371	LYS	Peptide
1	B	425	PHE	Peptide
1	C	129	ASN	Peptide
1	C	290	GLN	Peptide
1	C	371	LYS	Peptide
1	C	425	PHE	Peptide
1	D	290	GLN	Peptide
1	D	371	LYS	Peptide
1	D	425	PHE	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	4919	0	4975	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4978	0	5041	55	0
1	C	5217	0	5275	58	0
1	D	4919	0	4975	44	0
2	E	1165	0	1096	9	0
3	A	2	0	0	0	0
3	E	4	0	0	0	0
All	All	21204	0	21362	186	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 4.

All (186) 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:B:76:GLY:H	1:B:80:GLU:H	1.43	0.66
1:D:76:GLY:H	1:D:80:GLU:H	1.43	0.66
1:B:174:ASN:ND2	1:B:211:PHE:O	2.29	0.65
1:D:230:GLN:HE21	1:D:235:VAL:HG22	1.61	0.64
1:A:351:PRO:HB2	1:A:373:LEU:HB2	1.80	0.63
1:B:636:ARG:HH21	1:C:160:ILE:HG22	1.63	0.63
1:C:497:MET:HE1	1:D:568:LEU:HD21	1.80	0.62
1:D:280:LEU:HD22	1:D:631:LEU:HD23	1.82	0.61
1:A:174:ASN:ND2	1:A:211:PHE:O	2.33	0.61
2:E:27:ILE:HB	2:E:63:ILE:HB	1.83	0.60
1:C:75:ARG:NH2	1:C:111:THR:OG1	2.34	0.60
1:C:590:ASP:OD2	1:C:594:ARG:NH1	2.35	0.59
1:D:377:TYR:OH	1:D:443:ARG:NH2	2.36	0.59
1:D:564:ILE:HG13	1:D:568:LEU:HD23	1.83	0.59
1:D:128:GLN:NE2	2:E:38:SER:O	2.35	0.58
1:D:174:ASN:ND2	1:D:211:PHE:O	2.36	0.58
1:D:351:PRO:HB2	1:D:373:LEU:HB2	1.85	0.58
1:B:425:PHE:O	1:B:427:VAL:N	2.32	0.58
1:D:425:PHE:O	1:D:427:VAL:N	2.34	0.58
1:A:622:GLU:O	1:B:42:LYS:NZ	2.36	0.58
1:A:425:PHE:O	1:A:427:VAL:N	2.31	0.58
1:A:34:ASP:OD1	1:D:632:ARG:NH1	2.34	0.57
1:D:302:ARG:NH2	1:D:591:GLU:OE2	2.36	0.57
1:D:175:SER:HB3	1:D:178:ILE:HD13	1.86	0.57
1:C:425:PHE:O	1:C:427:VAL:N	2.34	0.57
1:A:619:CYS:SG	1:A:620:GLY:N	2.77	0.57
1:C:377:TYR:OH	1:C:443:ARG:NH2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:481:MET:HE1	1:A:582:HIS:HA	1.88	0.56
1:B:193:ASP:OD1	1:B:197:ASN:N	2.31	0.55
1:B:280:LEU:HD22	1:B:631:LEU:HD23	1.88	0.55
1:B:172:CYS:SG	1:B:206:GLN:NE2	2.79	0.55
1:C:564:ILE:HG13	1:C:568:LEU:HD23	1.89	0.55
1:B:486:ILE:HD13	1:C:569:LEU:HD22	1.88	0.54
1:B:351:PRO:HB2	1:B:373:LEU:HB2	1.89	0.54
1:A:193:ASP:OD1	1:A:197:ASN:N	2.37	0.54
1:B:257:GLN:NE2	1:B:310:GLN:OE1	2.41	0.54
1:B:511:ILE:HD11	1:B:554:MET:HB3	1.89	0.54
1:C:302:ARG:HH11	1:C:305:ARG:HH21	1.57	0.53
1:C:351:PRO:HB2	1:C:373:LEU:HB2	1.89	0.53
1:C:193:ASP:OD1	1:C:197:ASN:N	2.40	0.53
1:C:479:THR:HA	1:C:482:ILE:HG22	1.90	0.53
1:B:75:ARG:NH2	1:B:111:THR:OG1	2.42	0.53
1:A:516:ASP:HB3	1:D:367:LEU:HD21	1.91	0.53
1:D:193:ASP:OD1	1:D:197:ASN:N	2.39	0.53
2:E:47:GLU:O	2:E:51:MET:N	2.41	0.53
1:B:350:ARG:HH22	1:B:451:VAL:HG12	1.73	0.53
1:D:479:THR:HA	1:D:482:ILE:HG22	1.90	0.53
1:A:486:ILE:HD13	1:B:569:LEU:HD22	1.91	0.52
1:C:213:CYS:HA	1:C:255:MET:HE1	1.92	0.52
1:B:230:GLN:HE21	1:B:235:VAL:HG22	1.74	0.52
1:B:475:LEU:HD12	1:C:492:ARG:HB3	1.91	0.52
1:A:273:LEU:HD11	1:B:160:ILE:HG12	1.92	0.52
1:A:501:ILE:HD11	1:A:534:PHE:HA	1.92	0.52
1:B:525:ASP:N	1:B:525:ASP:OD1	2.43	0.52
1:C:643:ARG:HH22	2:E:133:ASP:HB3	1.74	0.52
1:B:75:ARG:NH1	1:B:108:GLU:OE2	2.43	0.52
1:B:58:VAL:O	1:B:62:ASN:N	2.40	0.51
1:A:459:VAL:HG13	1:A:460:LEU:HD12	1.92	0.51
2:E:26:THR:HB	2:E:62:THR:HB	1.93	0.51
1:B:640:ASN:HD22	2:E:90:ARG:HH21	1.59	0.50
1:D:236:PRO:HB2	1:D:240:GLY:HA2	1.94	0.50
1:B:190:ARG:NH1	1:B:227:ASP:OD2	2.39	0.50
2:E:92:PHE:HB3	2:E:100:ILE:HD13	1.94	0.50
1:C:632:ARG:NH1	1:D:34:ASP:OD1	2.40	0.49
1:B:479:THR:HA	1:B:482:ILE:HG22	1.93	0.49
1:B:481:MET:HE1	1:B:582:HIS:HA	1.93	0.49
1:A:568:LEU:HD21	1:D:497:MET:HE1	1.94	0.49
1:D:356:THR:OG1	1:D:357:ASN:N	2.40	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:ASP:OD1	1:A:196:GLY:N	2.45	0.49
1:A:479:THR:HA	1:A:482:ILE:HG22	1.94	0.49
1:A:237:ASN:OD1	1:A:241:LEU:N	2.46	0.49
1:C:284:ASP:O	1:C:291:SER:OG	2.26	0.49
1:D:130:MET:HB3	1:D:132:LEU:HB2	1.95	0.49
1:A:119:THR:HG23	1:A:121:LEU:H	1.78	0.48
1:A:478:PHE:HZ	1:B:577:MET:HE3	1.78	0.48
1:A:42:LYS:NZ	1:D:622:GLU:O	2.37	0.48
1:A:132:LEU:O	1:A:136:LEU:N	2.38	0.48
1:C:119:THR:HG23	1:C:121:LEU:H	1.79	0.48
1:C:175:SER:HB3	1:C:178:ILE:HD13	1.95	0.48
2:E:106:ARG:HG2	2:E:121:VAL:HG21	1.95	0.48
1:C:174:ASN:ND2	1:C:211:PHE:O	2.47	0.48
1:C:575:ILE:O	1:C:579:GLY:N	2.47	0.47
1:C:71:LYS:HB3	1:C:73:HIS:HD2	1.80	0.47
1:A:399:ILE:HD13	1:A:429:ILE:HD11	1.96	0.47
1:C:459:VAL:HG13	1:C:460:LEU:HD12	1.96	0.47
1:C:230:GLN:HE21	1:C:235:VAL:HG22	1.79	0.47
1:B:236:PRO:HB2	1:B:240:GLY:HA2	1.95	0.47
1:B:564:ILE:HG13	1:B:568:LEU:HD23	1.97	0.47
1:C:475:LEU:HD12	1:D:492:ARG:HB3	1.97	0.47
1:A:367:LEU:HD21	1:B:516:ASP:HB3	1.97	0.46
1:D:619:CYS:SG	1:D:620:GLY:N	2.88	0.46
1:C:356:THR:OG1	1:C:357:ASN:N	2.40	0.46
1:C:608:LEU:HD22	1:C:609:PRO:HD2	1.98	0.46
1:C:486:ILE:HD13	1:D:569:LEU:HD22	1.97	0.46
1:D:189:ILE:HD12	1:D:190:ARG:HG3	1.98	0.45
1:D:615:ARG:HD2	1:D:629:TRP:CE3	2.51	0.45
1:D:292:LEU:HD12	1:D:292:LEU:HA	1.79	0.45
1:C:132:LEU:HD12	1:C:135:ALA:HB3	1.98	0.45
1:C:276:THR:OG1	1:C:277:LEU:N	2.49	0.45
1:D:190:ARG:NH1	1:D:227:ASP:OD2	2.43	0.45
1:A:81:THR:OG1	1:A:84:HIS:ND1	2.39	0.45
1:A:564:ILE:HG13	1:A:568:LEU:HD23	1.99	0.45
1:B:356:THR:OG1	1:B:357:ASN:N	2.48	0.45
1:C:189:ILE:H	1:C:189:ILE:HG13	1.60	0.45
1:A:76:GLY:H	1:A:80:GLU:H	1.65	0.45
1:A:280:LEU:HD22	1:A:631:LEU:HD23	1.98	0.45
1:C:283:ILE:HG22	1:C:292:LEU:HD22	1.99	0.45
1:B:635:ASP:N	1:B:635:ASP:OD1	2.50	0.45
1:D:189:ILE:H	1:D:189:ILE:HG13	1.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:230:GLN:HB2	1:C:234:LEU:HD12	2.00	0.44
1:C:378:MET:SD	1:C:378:MET:N	2.77	0.44
1:A:45:TRP:HA	1:A:51:LEU:HD12	2.00	0.44
1:A:175:SER:HB3	1:A:178:ILE:HD13	1.99	0.44
1:C:45:TRP:HA	1:C:51:LEU:HD12	1.99	0.44
1:C:92:LEU:HD13	1:C:130:MET:HG3	1.99	0.44
1:C:619:CYS:SG	1:C:620:GLY:N	2.90	0.44
1:A:497:MET:HE2	1:A:537:PHE:HE2	1.83	0.43
1:B:190:ARG:HH22	1:B:227:ASP:HB2	1.83	0.43
1:A:119:THR:HG22	1:A:122:HIS:CE1	2.54	0.43
1:B:413:THR:HA	1:B:416:PHE:HD2	1.82	0.43
1:B:213:CYS:HA	1:B:255:MET:HE1	2.00	0.43
1:B:189:ILE:H	1:B:189:ILE:HG13	1.73	0.43
1:B:619:CYS:SG	1:B:620:GLY:N	2.91	0.43
1:C:413:THR:HA	1:C:416:PHE:HD2	1.83	0.43
1:C:694:ASN:O	1:C:698:LEU:N	2.44	0.43
1:D:112:SER:HA	1:D:116:GLU:HG2	2.00	0.43
1:B:119:THR:HG23	1:B:121:LEU:H	1.84	0.43
1:D:290:GLN:HB3	1:D:294:GLU:HB2	1.99	0.43
1:B:344:THR:HG22	1:B:458:LEU:HD11	2.01	0.43
1:D:608:LEU:HD22	1:D:609:PRO:HD2	2.00	0.43
1:C:75:ARG:NH1	1:C:108:GLU:OE2	2.52	0.43
1:B:497:MET:HE1	1:C:568:LEU:HD21	2.01	0.42
1:D:297:ILE:HD12	1:D:298:THR:HG23	2.01	0.42
1:A:525:ASP:OD1	1:A:525:ASP:N	2.51	0.42
1:B:193:ASP:OD1	1:B:196:GLY:N	2.52	0.42
1:C:522:HIS:CD2	1:C:544:PRO:HB3	2.54	0.42
1:C:236:PRO:HB2	1:C:240:GLY:HA2	2.01	0.42
2:E:129:ASP:OD2	2:E:134:GLY:N	2.50	0.42
1:A:230:GLN:HG3	1:A:234:LEU:HB2	2.00	0.42
1:C:504:PHE:HD1	1:C:504:PHE:HA	1.75	0.42
1:A:213:CYS:HA	1:A:255:MET:HE1	2.00	0.42
1:B:124:ALA:HB2	1:B:132:LEU:HD23	2.01	0.42
1:B:636:ARG:HG2	1:B:638:ASP:H	1.84	0.42
1:C:160:ILE:HD13	1:C:162:PHE:HD2	1.84	0.42
1:D:486:ILE:HD12	1:D:486:ILE:HA	1.89	0.42
1:D:522:HIS:CD2	1:D:544:PRO:HB3	2.55	0.42
1:B:408:PHE:HD1	1:B:408:PHE:HA	1.77	0.42
1:C:129:ASN:OD1	1:C:131:ASN:N	2.40	0.42
1:C:481:MET:HE1	1:C:582:HIS:HA	2.02	0.42
1:C:497:MET:HE2	1:C:537:PHE:HE2	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:538:LEU:HD21	1:D:568:LEU:HD22	2.01	0.42
1:D:625:LEU:HD22	1:D:625:LEU:HA	1.91	0.42
1:A:362:PRO:HD2	1:A:366:THR:HG21	2.02	0.41
1:A:632:ARG:NH1	1:B:34:ASP:OD1	2.41	0.41
1:D:137:LEU:HD23	1:D:137:LEU:HA	1.89	0.41
1:A:395:GLY:HA2	1:A:398:ILE:HD12	2.01	0.41
1:A:618:ILE:N	1:A:630:PHE:O	2.48	0.41
1:A:369:GLN:H	1:B:514:THR:HG22	1.84	0.41
1:A:373:LEU:HD13	1:A:376:ALA:HB3	2.01	0.41
1:A:344:THR:HG22	1:A:458:LEU:HD11	2.01	0.41
1:B:565:ILE:HD13	1:B:565:ILE:HA	1.92	0.41
1:C:253:THR:O	1:C:257:GLN:N	2.50	0.41
1:D:284:ASP:O	1:D:291:SER:OG	2.30	0.41
1:A:160:ILE:HG12	1:D:273:LEU:HD11	2.02	0.41
1:A:283:ILE:HG22	1:A:292:LEU:HD22	2.01	0.41
1:B:283:ILE:HG22	1:B:292:LEU:HD22	2.02	0.41
1:C:54:LYS:HE3	1:C:89:TYR:CZ	2.56	0.41
1:A:639:LEU:HD13	1:A:640:ASN:H	1.85	0.41
1:B:136:LEU:HD23	1:B:136:LEU:HA	1.93	0.41
1:B:610:ARG:HB3	1:B:611:CYS:H	1.73	0.41
1:B:636:ARG:NH1	1:C:207:PRO:HG3	2.36	0.41
1:D:520:LEU:HD23	1:D:520:LEU:HA	1.94	0.41
1:C:356:THR:HG1	1:C:357:ASN:H	1.65	0.41
1:B:353:LYS:HD2	1:B:373:LEU:HD23	2.04	0.40
1:C:189:ILE:HD12	1:C:190:ARG:HG3	2.02	0.40
1:C:625:LEU:HD22	1:C:625:LEU:HA	1.90	0.40
1:B:175:SER:HB3	1:B:178:ILE:HD13	2.03	0.40
1:B:358:ASN:OD1	1:B:358:ASN:N	2.52	0.40
1:D:452:VAL:HG13	1:D:453:PRO:HD3	2.03	0.40
1:A:378:MET:SD	1:A:378:MET:N	2.81	0.40
1:C:193:ASP:OD1	1:C:196:GLY:N	2.54	0.40
1:B:367:LEU:HD21	1:C:516:ASP:HB3	2.04	0.40
1:C:444:LEU:HD23	1:C:444:LEU:HA	1.96	0.40
1:D:192:GLN:HB3	1:D:196:GLY:HA2	2.03	0.40
1:A:522:HIS:CD2	1:A:544:PRO:HB3	2.56	0.40
1:B:616:SER:OG	1:B:617:GLY:N	2.54	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	611/725 (84%)	541 (88%)	67 (11%)	3 (0%)	24	59
1	B	617/725 (85%)	542 (88%)	73 (12%)	2 (0%)	36	69
1	C	643/725 (89%)	580 (90%)	61 (10%)	2 (0%)	36	69
1	D	611/725 (84%)	544 (89%)	64 (10%)	3 (0%)	24	59
2	E	146/149 (98%)	132 (90%)	14 (10%)	0	100	100
All	All	2628/3049 (86%)	2339 (89%)	279 (11%)	10 (0%)	31	64

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	426	HIS
1	B	426	HIS
1	C	426	HIS
1	D	426	HIS
1	C	372	LEU
1	A	473	GLN
1	B	372	LEU
1	D	372	LEU
1	A	372	LEU
1	D	621	ARG

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	532/633 (84%)	524 (98%)	8 (2%)	57	70
1	B	538/633 (85%)	528 (98%)	10 (2%)	50	66
1	C	562/633 (89%)	557 (99%)	5 (1%)	70	75
1	D	532/633 (84%)	524 (98%)	8 (2%)	57	70
2	E	126/127 (99%)	125 (99%)	1 (1%)	73	77
All	All	2290/2659 (86%)	2258 (99%)	32 (1%)	57	71

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	ILE
1	A	183	ILE
1	A	189	ILE
1	A	371	LYS
1	A	551	LEU
1	A	608	LEU
1	A	625	LEU
1	A	639	LEU
1	B	160	ILE
1	B	181	LEU
1	B	183	ILE
1	B	189	ILE
1	B	377	TYR
1	B	459	VAL
1	B	492	ARG
1	B	551	LEU
1	B	608	LEU
1	B	625	LEU
1	C	181	LEU
1	C	189	ILE
1	C	365	ASN
1	C	551	LEU
1	C	625	LEU
1	D	160	ILE
1	D	181	LEU
1	D	189	ILE
1	D	459	VAL
1	D	492	ARG
1	D	551	LEU
1	D	608	LEU
1	D	625	LEU

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Mol	Chain	Res	Type
2	E	9	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	165	HIS
1	A	206	GLN
1	A	257	GLN
1	A	365	ASN
1	A	370	GLN
1	B	158	ASN
1	B	165	HIS
1	B	197	ASN
1	B	206	GLN
1	B	257	GLN
1	B	258	HIS
1	B	365	ASN
1	B	370	GLN
1	B	640	ASN
1	C	73	HIS
1	C	257	GLN
1	C	365	ASN
1	C	587	HIS
1	D	40	GLN
1	D	165	HIS
1	D	230	GLN
1	D	257	GLN
1	D	306	GLN
1	D	365	ASN
1	D	370	GLN
2	E	53	ASN
2	E	107	HIS
2	E	111	ASN

5.3.3 RNA ⓘ

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 6 ligands modelled in this entry, 6 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.

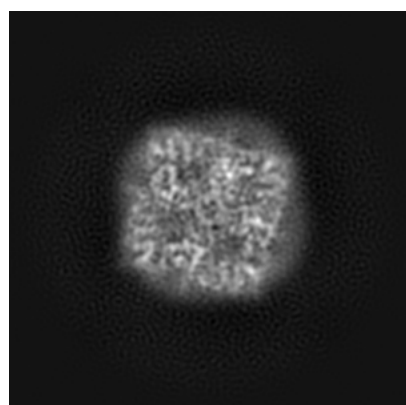
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8961. These allow visual inspection of the internal detail of the map and identification of artifacts.

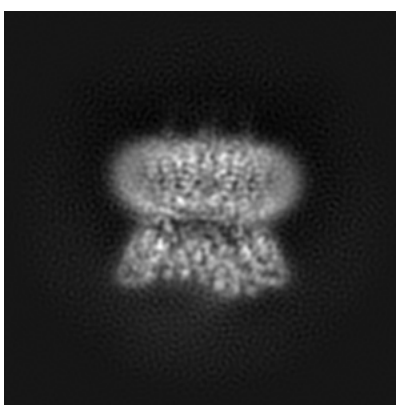
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

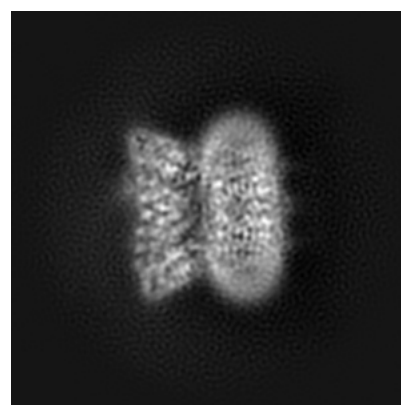
6.1.1 Primary map



X



Y

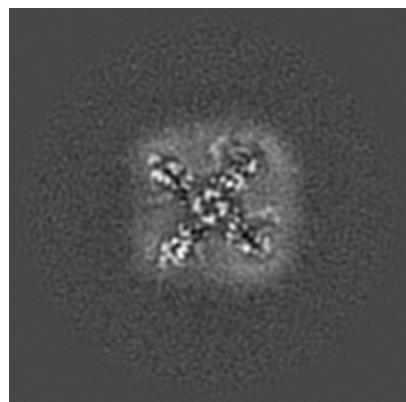


Z

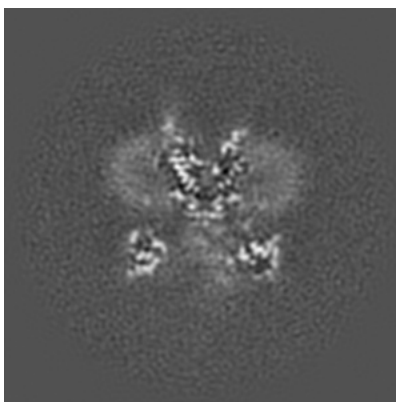
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

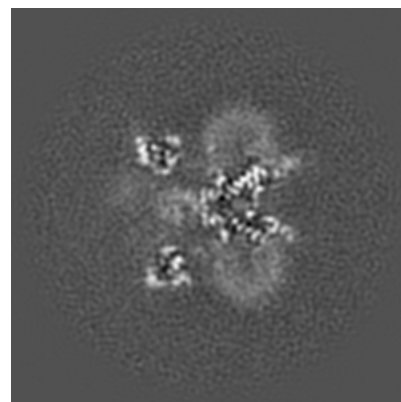
6.2.1 Primary map



X Index: 125



Y Index: 125

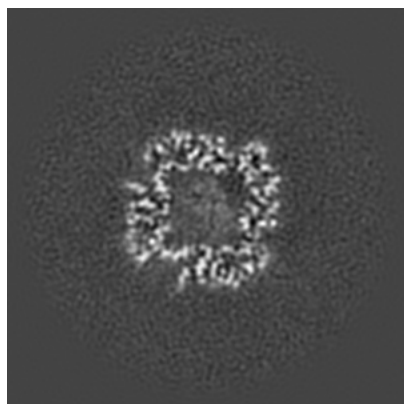


Z Index: 125

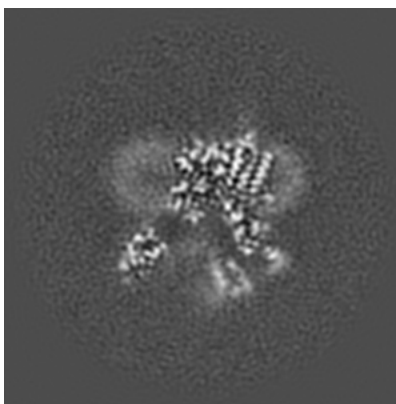
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

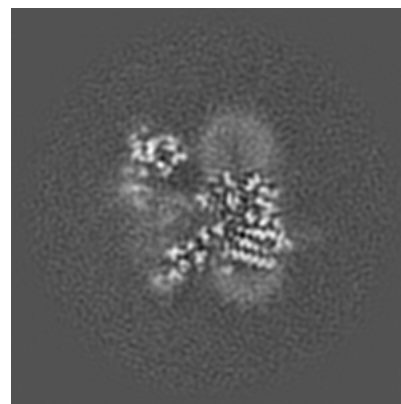
6.3.1 Primary map



X Index: 101



Y Index: 137

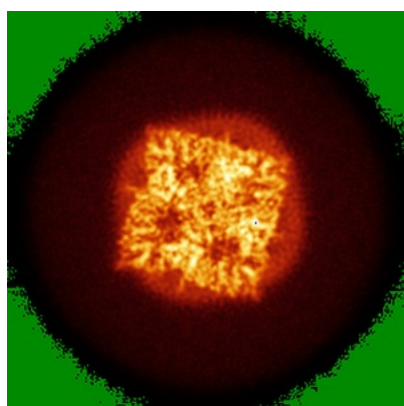


Z Index: 136

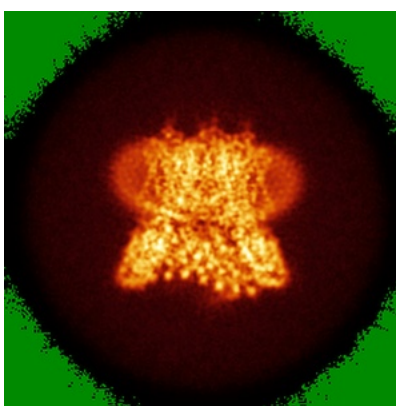
The images above show the largest variance slices of the map in three orthogonal directions.

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

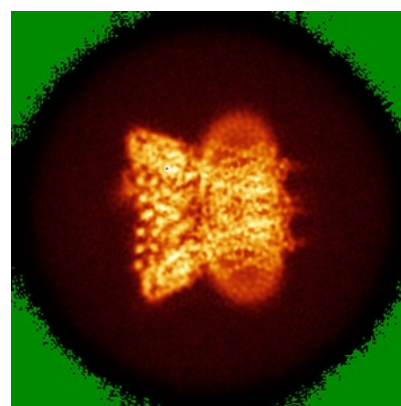
6.4.1 Primary map



X



Y

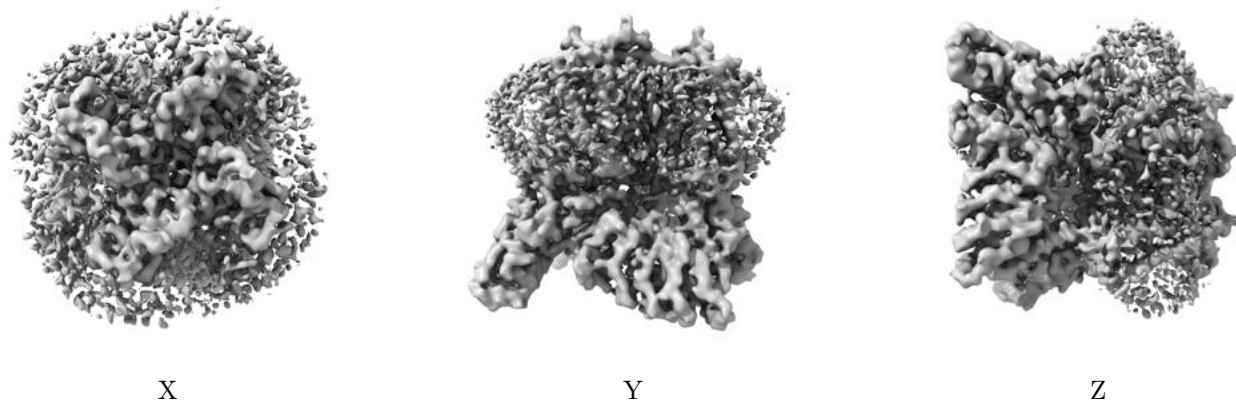


Z

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

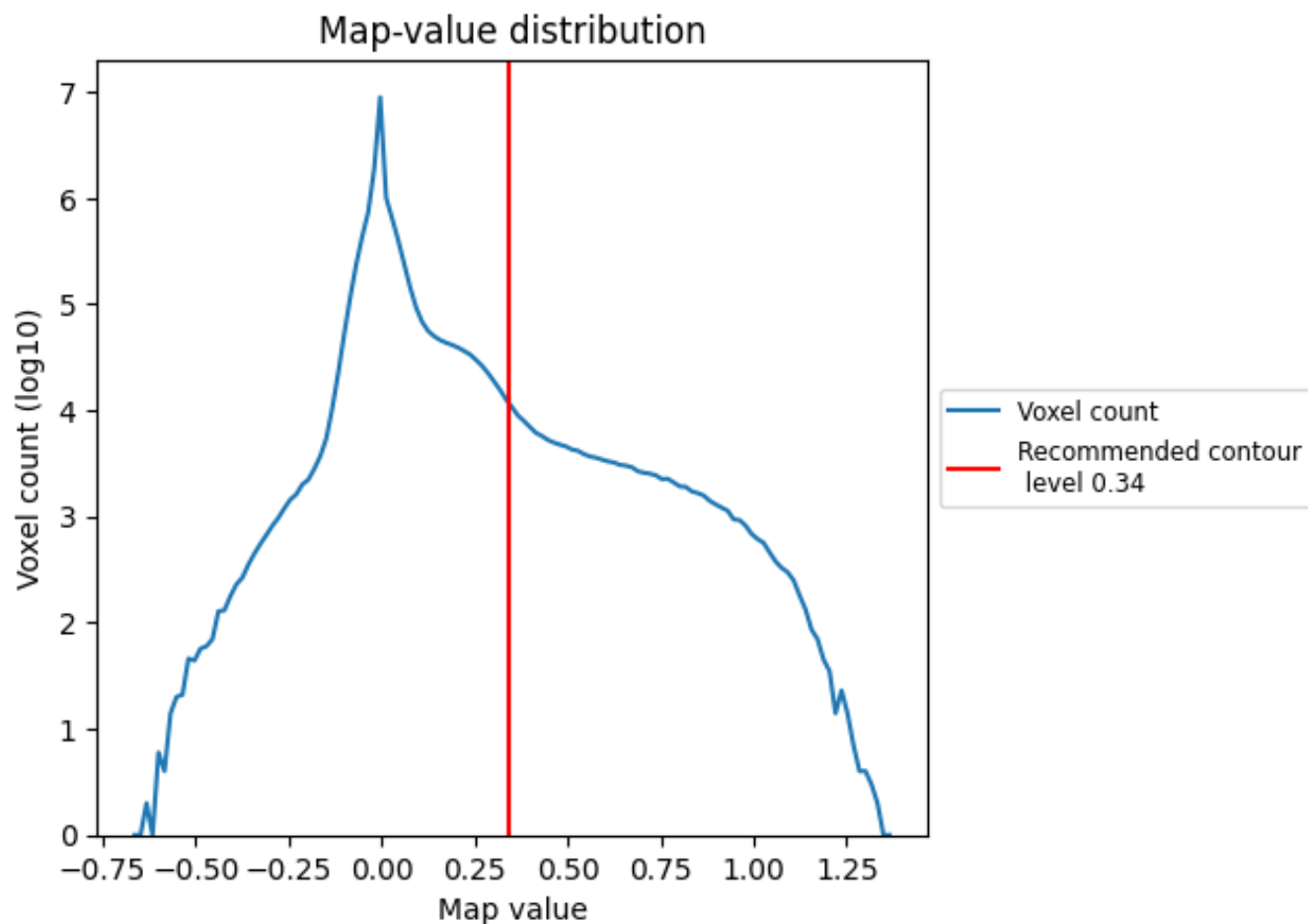
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

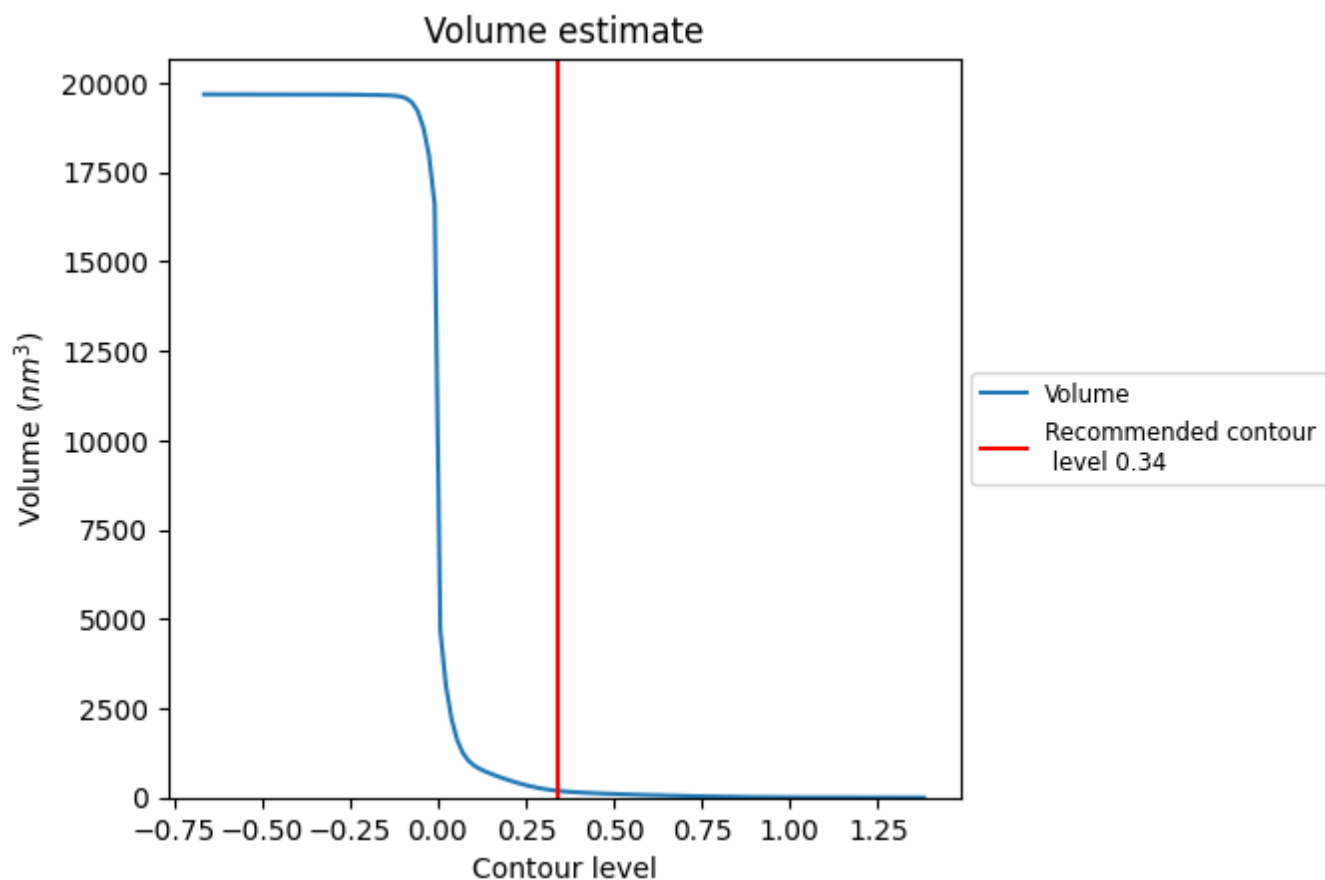
This section contains the results of statistical analysis of the map.

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.

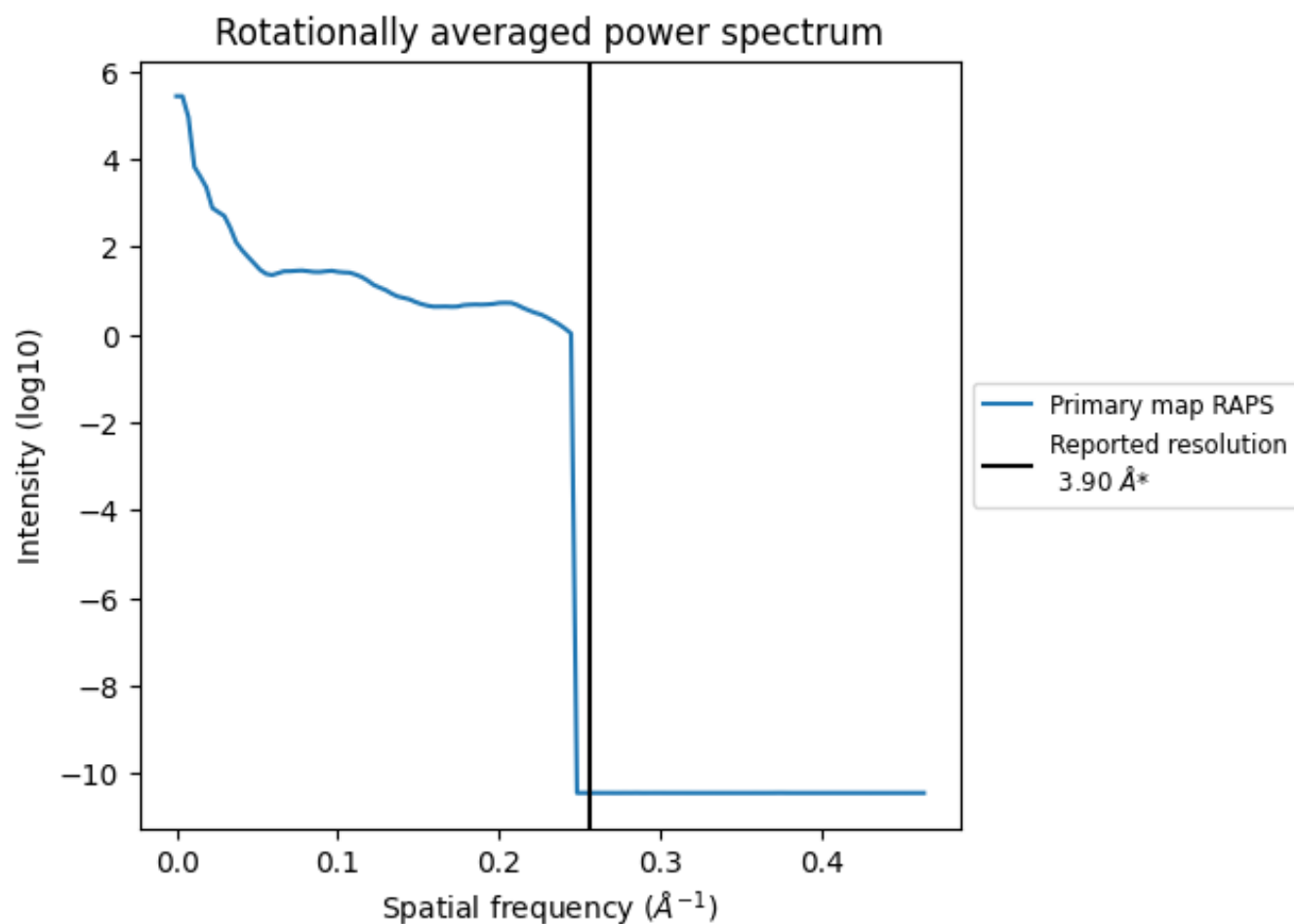
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 190 nm³; this corresponds to an approximate mass of 171 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 0.256 Å⁻¹

8 Fourier-Shell correlation

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

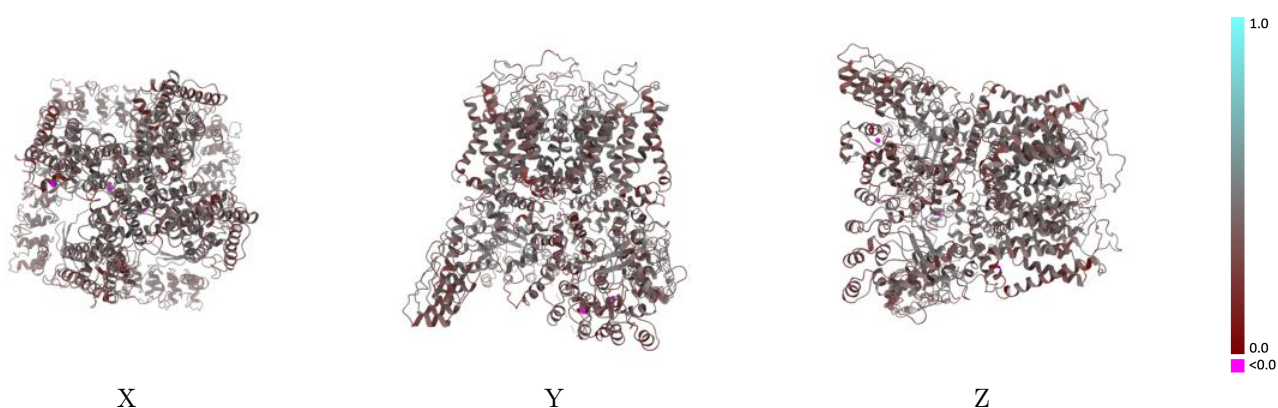
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8961 and PDB model 6E2F. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)

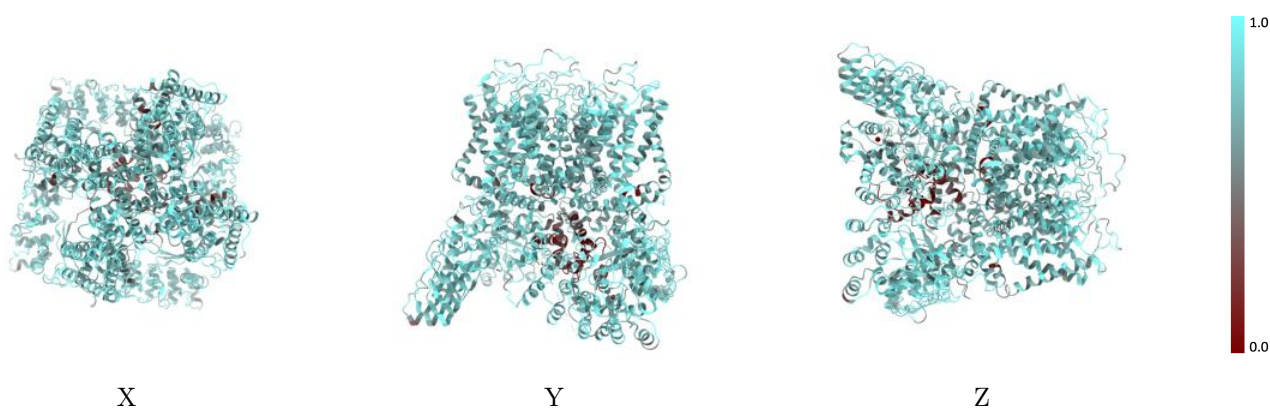
This section was not generated.

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



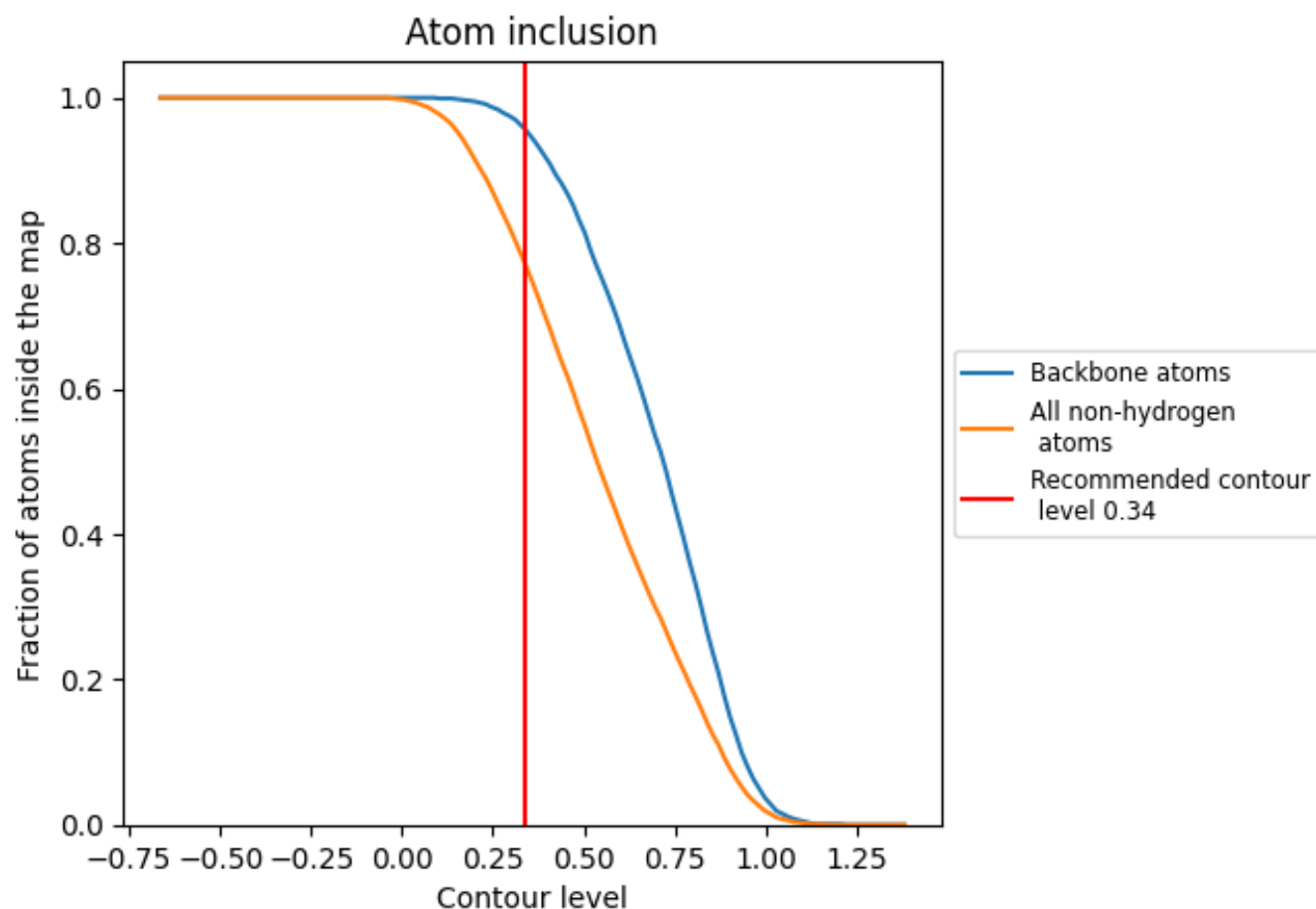
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.34).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% 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 (0.34) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7700	0.3920
A	0.7910	0.3980
B	0.7840	0.3930
C	0.7790	0.3960
D	0.7950	0.3980
E	0.4830	0.3100

