



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 01:35 AM UTC

PDB ID : 6E2G / pdb_00006e2g
EMDB ID : EMD-8962
Title : Cryo-EM structure of rat TRPV6 in complex with Calmodulin
Authors : Singh, A.K.; McGoldrick, L.L.; Sobolevsky, A.I.
Deposited on : 2018-07-11
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:

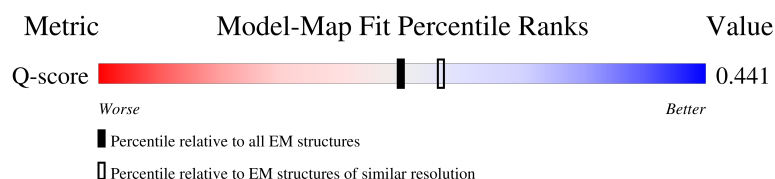
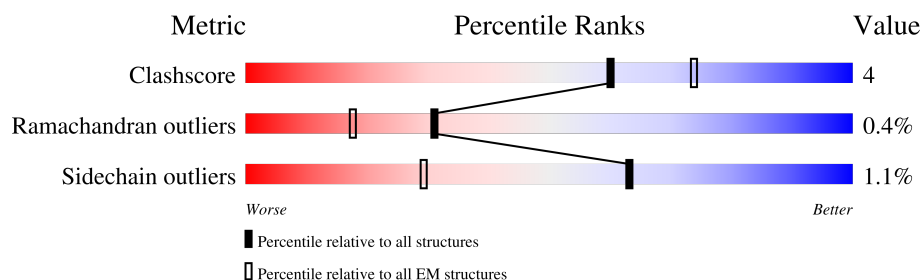
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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	727	 7% 72% 12% • 16%
1	B	727	 8% 74% 11% • 15%
1	C	727	 10% 77% 11% 11%
1	D	727	 7% 72% 12% 16%

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Mol	Chain	Length	Quality of chain
2	E	149	63% 81% 18%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 21108 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	611	Total	C	N	O	S	0	0
			4885	3166	818	866	35		
1	B	621	Total	C	N	O	S	0	0
			4979	3223	842	879	35		
1	C	644	Total	C	N	O	S	0	0
			5172	3337	884	916	35		
1	D	613	Total	C	N	O	S	0	0
			4901	3176	821	869	35		

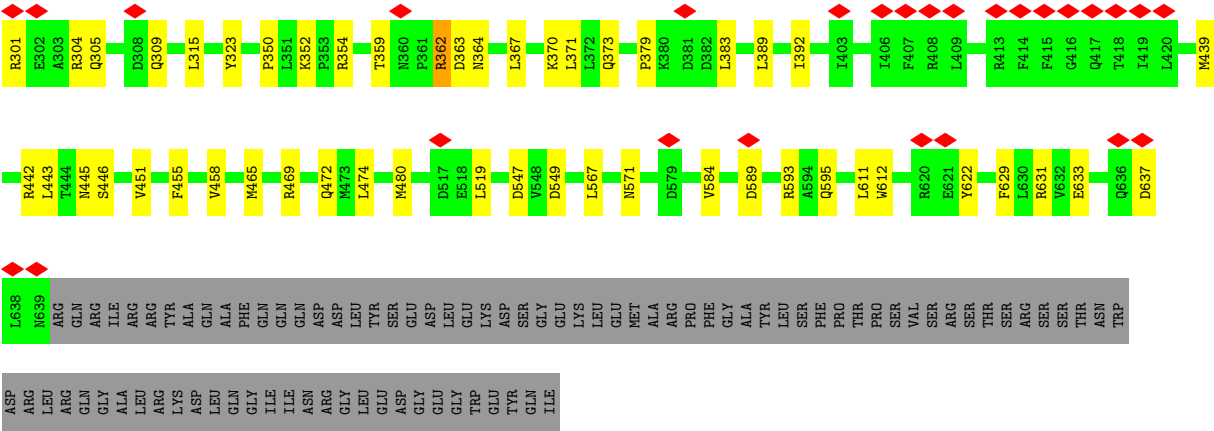
- Molecule 2 is a protein called Calmodulin-1.

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

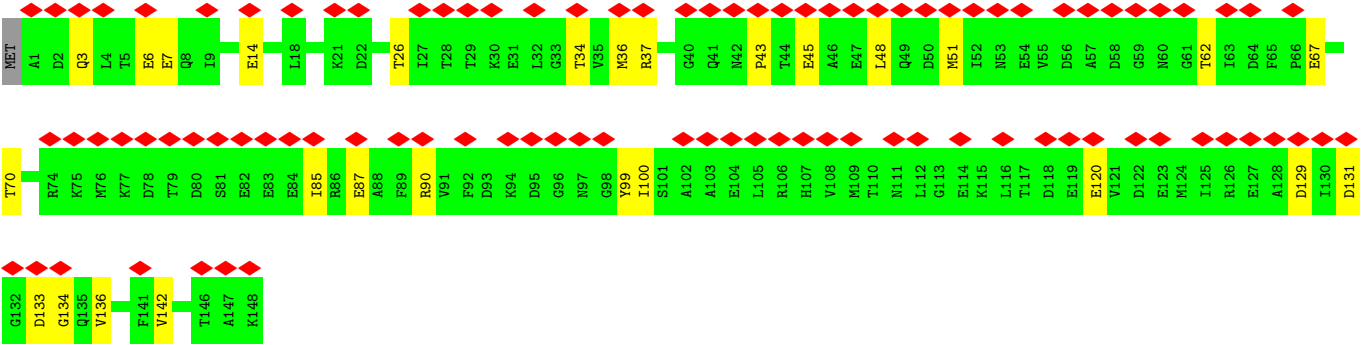
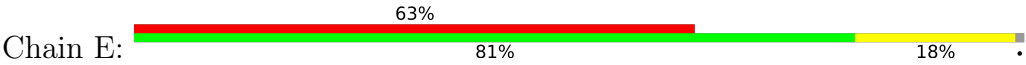
- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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





● Molecule 2: Calmodulin-1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	111593	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 ($e^-/\text{\AA}^2$)	47.0	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	1.646	Depositor
Minimum map value	-0.963	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.078	Depositor
Recommended contour level	0.4	Depositor
Map size (Å)	254.4, 254.4, 254.4	wwPDB
Map dimensions	240, 240, 240	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

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	0.43	0/4998	0.67	1/6787 (0.0%)
1	B	0.43	0/5093	0.68	3/6913 (0.0%)
1	C	0.44	0/5288	0.68	3/7173 (0.0%)
1	D	0.43	0/5014	0.69	1/6809 (0.0%)
2	E	0.25	0/1177	0.53	0/1580
All	All	0.43	0/21570	0.67	8/29262 (0.0%)

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	4
1	C	0	2
1	D	0	2
All	All	0	11

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	637	ASP	CA-C-N	6.94	134.79	121.54
1	B	637	ASP	C-N-CA	6.94	134.79	121.54
1	C	209	LYS	CA-C-N	5.56	132.15	121.54
1	C	209	LYS	C-N-CA	5.56	132.15	121.54
1	A	211	PHE	N-CA-C	-5.25	108.13	114.75
1	D	211	PHE	N-CA-C	-5.17	108.24	114.75
1	B	310	THR	N-CA-C	5.12	121.13	109.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	310	THR	N-CA-C	5.02	120.89	109.81

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	296	VAL	Peptide
1	A	358	ARG	Peptide
1	A	359	THR	Peptide
1	B	229	LYS	Peptide
1	B	286	GLY	Peptide
1	B	362	ARG	Peptide
1	B	585	ALA	Peptide
1	C	209	LYS	Peptide
1	C	362	ARG	Peptide
1	D	362	ARG	Peptide
1	D	71	GLU	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	4885	0	4937	51	0
1	B	4979	0	5039	50	0
1	C	5172	0	5220	47	0
1	D	4901	0	4954	52	0
2	E	1165	0	1096	19	0
3	A	2	0	0	0	0
3	E	4	0	0	0	0
All	All	21108	0	21246	190	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 (190) 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:304:ARG:HH12	1:A:420:LEU:HD11	1.58	0.68
1:B:151:VAL:O	1:B:158:ASN:ND2	2.27	0.68
1:B:469:ARG:NH2	1:B:595:GLN:OE1	2.28	0.66
1:A:500:ILE:HD11	1:A:533:PHE:HA	1.78	0.65
1:D:567:LEU:HA	1:D:571:ASN:HD22	1.62	0.64
1:D:469:ARG:NH2	1:D:595:GLN:OE1	2.33	0.61
1:B:359:THR:HB	1:B:364:ASN:HB3	1.82	0.60
1:A:118:GLN:NE2	1:D:269:TYR:OH	2.34	0.60
1:C:256:GLN:NE2	1:C:309:GLN:OE1	2.34	0.60
1:C:354:ARG:HD2	1:C:367:LEU:HA	1.82	0.60
1:A:34:ASP:OD1	1:D:631:ARG:NH1	2.30	0.59
1:D:190:ARG:NH2	1:D:229:LYS:O	2.35	0.59
1:C:469:ARG:NH2	1:C:595:GLN:OE1	2.35	0.59
1:C:631:ARG:HH12	1:D:37:ASN:HD22	1.51	0.58
1:C:500:ILE:HD11	1:C:533:PHE:HA	1.83	0.58
1:A:256:GLN:NE2	1:A:309:GLN:OE1	2.35	0.58
1:C:151:VAL:O	1:C:158:ASN:ND2	2.36	0.58
1:D:373:GLN:H	1:D:445:ASN:HD21	1.51	0.58
1:B:619:GLY:HA2	1:B:624:LEU:HD23	1.86	0.57
1:D:354:ARG:HD2	1:D:367:LEU:HA	1.85	0.57
1:A:348:TYR:O	1:A:384:ARG:NH1	2.35	0.57
1:B:279:LEU:HD11	1:B:315:LEU:HD21	1.86	0.57
1:D:589:ASP:O	1:D:593:ARG:NH1	2.38	0.57
1:D:90:ASP:OD2	2:E:37:ARG:NH2	2.37	0.56
1:D:370:LYS:NZ	1:D:446:SER:O	2.38	0.56
1:A:567:LEU:HA	1:A:571:ASN:HD22	1.71	0.56
1:D:114:LEU:HA	1:D:150:SER:HB3	1.88	0.56
1:A:268:THR:HG22	1:A:273:THR:HG22	1.88	0.56
1:A:547:ASP:O	1:D:362:ARG:NH1	2.39	0.56
1:C:620:ARG:NH2	1:C:626:ASP:OD1	2.39	0.56
1:A:474:LEU:HD11	1:B:495:LEU:HD23	1.87	0.55
1:A:42:LYS:HD2	1:D:622:TYR:HA	1.89	0.55
1:A:279:LEU:HD11	1:A:315:LEU:HD21	1.89	0.55
1:C:699:ARG:O	1:C:703:GLY:N	2.40	0.55
1:B:362:ARG:NH1	1:C:547:ASP:O	2.40	0.55
1:B:54:LYS:O	1:B:91:ASN:ND2	2.40	0.55
1:A:54:LYS:O	1:A:91:ASN:ND2	2.40	0.54
1:C:511:PHE:HD1	1:C:514:GLU:HG3	1.73	0.54
1:B:183:ILE:HD13	1:B:187:ALA:HB3	1.89	0.54
1:A:360:ASN:H	1:A:364:ASN:HD22	1.54	0.53
1:A:537:LEU:HD21	1:B:567:LEU:HD23	1.89	0.53
1:D:193:ASP:OD1	1:D:197:ASN:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:GLN:NE2	1:A:78:MET:SD	2.75	0.53
1:C:42:LYS:NZ	1:C:46:GLU:OE2	2.40	0.53
1:D:183:ILE:HD13	1:D:187:ALA:HB3	1.91	0.53
1:B:193:ASP:OD1	1:B:197:ASN:N	2.42	0.53
1:B:256:GLN:NE2	1:B:309:GLN:OE1	2.42	0.53
1:B:44:ILE:HG13	1:B:50:LEU:HB3	1.91	0.53
1:B:80:GLU:HG3	1:B:85:ILE:HD11	1.91	0.52
1:A:208:ASN:ND2	1:D:637:ASP:OD2	2.42	0.52
2:E:7:GLU:OE2	2:E:90:ARG:NE	2.43	0.52
1:A:413:ARG:NH2	1:A:420:LEU:O	2.42	0.52
1:D:252:ILE:HD11	1:D:305:GLN:HG3	1.92	0.52
1:B:563:ILE:HG23	1:B:567:LEU:HD11	1.92	0.52
1:C:296:VAL:HG12	1:C:420:LEU:HD13	1.91	0.52
1:A:256:GLN:O	1:A:260:GLN:NE2	2.43	0.52
1:B:116:GLU:HB3	1:B:149:GLY:HA2	1.92	0.52
1:C:224:GLY:HA2	1:C:231:LEU:HD22	1.92	0.52
1:D:350:PRO:HB2	1:D:371:LEU:HB2	1.91	0.52
1:A:360:ASN:HB3	1:A:364:ASN:HB2	1.90	0.52
1:D:276:LEU:HD12	1:D:629:PHE:HB3	1.92	0.52
1:D:352:LYS:HG3	1:D:371:LEU:HD22	1.92	0.52
1:D:256:GLN:HE21	1:D:309:GLN:HE21	1.58	0.51
1:A:183:ILE:HD13	1:A:187:ALA:HB3	1.92	0.51
1:C:692:ARG:NH2	2:E:120:GLU:OE2	2.44	0.51
1:A:269:TYR:OH	1:B:118:GLN:NE2	2.43	0.51
1:D:480:MET:HE1	1:D:584:VAL:HG23	1.91	0.51
1:B:620:ARG:NH2	1:B:626:ASP:OD1	2.34	0.51
1:C:199:VAL:HA	1:C:202:ILE:HD12	1.93	0.51
1:B:279:LEU:HD22	1:B:630:LEU:HB2	1.93	0.50
1:C:642:ARG:HH22	2:E:133:ASP:HB3	1.76	0.50
1:B:318:LEU:HB3	1:B:322:ARG:HH21	1.76	0.50
2:E:100:ILE:HG23	2:E:136:VAL:HB	1.94	0.50
1:B:167:LEU:HD21	1:B:189:ILE:HD13	1.93	0.49
1:A:469:ARG:NH2	1:A:595:GLN:OE1	2.45	0.49
1:A:495:LEU:HD23	1:D:474:LEU:HD11	1.94	0.49
1:D:42:LYS:NZ	1:D:46:GLU:OE2	2.44	0.49
1:C:639:ASN:ND2	2:E:131:ASP:O	2.42	0.49
1:D:389:LEU:HA	1:D:392:ILE:HG22	1.94	0.49
1:A:33:ARG:O	1:A:37:ASN:ND2	2.46	0.49
1:C:376:TYR:HE1	1:C:384:ARG:HH21	1.61	0.49
1:A:619:GLY:HA2	1:A:624:LEU:HD23	1.95	0.49
1:B:348:TYR:O	1:B:384:ARG:NH1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:ARG:HD2	1:B:367:LEU:HA	1.95	0.49
1:D:279:LEU:O	1:D:284:SER:OG	2.30	0.48
1:B:640:ARG:HH21	2:E:87:GLU:HB2	1.77	0.48
1:C:352:LYS:HG3	1:C:371:LEU:HD22	1.95	0.48
1:C:362:ARG:NH1	1:D:547:ASP:O	2.46	0.48
1:A:209:LYS:HD3	1:A:251:ASN:HD21	1.77	0.48
1:B:352:LYS:N	1:B:369:GLN:O	2.46	0.48
1:A:466:TYR:HD1	1:A:469:ARG:HH11	1.61	0.48
1:C:300:LYS:NZ	1:C:302:GLU:OE2	2.38	0.48
1:D:172:CYS:SG	1:D:173:VAL:N	2.86	0.48
1:A:167:LEU:HD21	1:A:189:ILE:HD13	1.94	0.48
1:D:279:LEU:HD12	1:D:282:ILE:HB	1.95	0.48
1:A:511:PHE:HD1	1:A:514:GLU:HG3	1.78	0.48
2:E:26:THR:HB	2:E:62:THR:HB	1.94	0.48
1:B:354:ARG:NH2	1:B:366:LEU:O	2.44	0.48
1:D:350:PRO:HD2	1:D:442:ARG:HH22	1.78	0.47
1:C:651:GLN:HA	2:E:51:MET:HE3	1.96	0.47
1:B:49:LEU:O	1:B:53:ALA:N	2.48	0.47
1:D:161:TYR:OH	1:D:193:ASP:OD2	2.32	0.47
1:B:479:ILE:HD11	1:B:595:GLN:HG3	1.96	0.46
1:C:209:LYS:O	1:C:211:PHE:N	2.35	0.46
1:C:534:GLU:HG2	1:C:539:ILE:HD11	1.98	0.46
1:D:359:THR:HG23	1:D:364:ASN:HB2	1.97	0.46
1:C:484:MET:HE1	1:C:580:THR:HG23	1.98	0.46
1:B:268:THR:HG22	1:B:273:THR:HG22	1.97	0.46
1:A:45:TRP:HA	1:A:51:LEU:HD12	1.98	0.45
1:B:269:TYR:OH	1:C:118:GLN:NE2	2.49	0.45
2:E:34:THR:O	2:E:37:ARG:NH1	2.48	0.45
1:C:62:ILE:HG23	1:C:66:LYS:HG3	1.98	0.45
1:A:146:ARG:NH1	1:A:164:GLU:OE1	2.50	0.45
1:C:112:SER:OG	1:C:113:GLU:N	2.48	0.45
1:D:33:ARG:HG2	1:D:114:LEU:HD21	1.98	0.45
1:C:339:ILE:O	1:C:343:THR:N	2.43	0.45
1:C:509:ILE:HD13	1:C:509:ILE:HA	1.84	0.45
1:A:84:HIS:O	1:A:88:LEU:N	2.43	0.45
1:B:212:ALA:O	1:B:216:TYR:N	2.50	0.45
1:B:631:ARG:HH12	1:C:37:ASN:HD22	1.63	0.45
1:C:267:TRP:HB2	1:D:88:LEU:HD21	1.99	0.45
1:A:497:ALA:HA	1:A:500:ILE:HG22	1.98	0.45
1:A:44:ILE:HD13	1:A:54:LYS:HE3	1.98	0.45
1:A:591:LEU:O	1:A:595:GLN:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:362:ARG:NH2	1:D:363:ASP:OD2	2.50	0.45
1:D:209:LYS:O	1:D:211:PHE:N	2.44	0.44
1:D:301:ARG:HH11	1:D:304:ARG:HH22	1.64	0.44
2:E:14:GLU:HG3	2:E:99:TYR:HE2	1.83	0.44
1:D:112:SER:OG	1:D:113:GLU:N	2.51	0.44
1:C:212:ALA:O	1:C:216:TYR:N	2.42	0.44
1:D:455:PHE:HA	1:D:458:VAL:HG22	2.00	0.44
1:A:458:VAL:HG23	1:A:459:LEU:HD12	2.00	0.44
1:C:350:PRO:HB2	1:C:371:LEU:HD23	2.00	0.44
1:C:643:ILE:H	1:C:643:ILE:HG13	1.56	0.44
2:E:67:GLU:HA	2:E:70:THR:HG22	1.99	0.44
1:C:350:PRO:HD2	1:C:442:ARG:HH22	1.83	0.43
1:C:607:LEU:HD23	1:C:611:LEU:HD11	1.99	0.43
1:A:75:LYS:HB3	1:A:79:GLY:HA2	2.01	0.43
1:B:40:GLN:NE2	1:B:78:MET:SD	2.91	0.43
1:B:283:ASP:N	1:B:283:ASP:OD1	2.51	0.43
1:C:363:ASP:HB2	1:D:549:ASP:HB2	2.00	0.43
1:D:472:GLN:H	1:D:472:GLN:HG2	1.47	0.43
1:A:489:LEU:HD12	1:A:489:LEU:HA	1.78	0.43
1:B:614:ARG:HG2	1:B:628:TRP:CE2	2.54	0.43
1:B:539:ILE:HG22	1:C:538:THR:HG23	2.00	0.43
1:B:315:LEU:O	1:B:319:LYS:N	2.40	0.43
1:B:534:GLU:HG2	1:B:539:ILE:HD11	2.01	0.43
1:D:249:GLU:OE1	1:D:300:LYS:NZ	2.52	0.43
1:D:279:LEU:HD11	1:D:315:LEU:HD21	2.00	0.43
1:A:611:LEU:H	1:A:611:LEU:HG	1.69	0.42
1:A:350:PRO:HB2	1:A:371:LEU:HD23	2.01	0.42
1:A:354:ARG:NH2	1:A:366:LEU:O	2.53	0.42
1:D:196:GLY:HA3	1:D:237:ASN:HB2	2.01	0.42
1:C:75:LYS:HB3	1:C:79:GLY:HA2	2.00	0.42
1:C:235:PRO:HB2	1:C:239:GLY:HA2	2.01	0.42
1:A:527:MET:HE2	1:B:546:TYR:HE1	1.85	0.42
1:A:199:VAL:HG13	1:A:200:LEU:HD22	2.01	0.42
1:B:631:ARG:HH12	1:C:37:ASN:ND2	2.18	0.42
1:C:192:GLN:HB3	1:C:196:GLY:HA2	2.01	0.42
1:D:78:MET:SD	1:D:78:MET:N	2.93	0.42
2:E:85:ILE:HG21	2:E:142:VAL:HG23	2.02	0.41
1:A:376:TYR:HE1	1:A:384:ARG:HH21	1.69	0.41
1:B:643:ILE:O	2:E:3:GLN:NE2	2.53	0.41
1:D:323:TYR:HB3	1:D:612:TRP:NE1	2.35	0.41
2:E:36:MET:HB3	2:E:43:PRO:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:ILE:O	1:B:290:SER:OG	2.38	0.41
1:C:370:LYS:NZ	1:C:446:SER:O	2.42	0.41
1:D:36:GLN:O	1:D:40:GLN:N	2.46	0.41
1:B:465:MET:HE2	1:B:465:MET:HB2	1.89	0.41
1:B:530:PHE:HE1	1:C:559:ALA:HB2	1.85	0.41
1:D:439:MET:HE3	1:D:443:LEU:HD21	2.03	0.41
1:A:536:PHE:HZ	1:A:569:MET:HG2	1.86	0.41
2:E:129:ASP:OD2	2:E:133:ASP:N	2.54	0.41
1:B:451:VAL:HB	1:C:553:MET:HE3	2.02	0.41
1:D:379:PRO:O	1:D:383:LEU:N	2.52	0.41
1:D:631:ARG:NH2	1:D:633:GLU:OE1	2.54	0.41
2:E:45:GLU:HA	2:E:48:LEU:HD12	2.03	0.41
1:A:363:ASP:HB2	1:B:549:ASP:HB2	2.03	0.41
1:A:398:ILE:HA	1:A:401:VAL:HG12	2.03	0.41
1:A:465:MET:HE2	1:A:465:MET:HB2	1.93	0.41
1:B:519:LEU:HG	1:B:550:LEU:HD11	2.03	0.41
1:D:465:MET:HE2	1:D:465:MET:HB2	1.86	0.41
1:A:200:LEU:HD13	1:A:200:LEU:HA	1.80	0.41
1:B:229:LYS:HA	1:B:229:LYS:HD2	1.85	0.40
2:E:129:ASP:OD2	2:E:134:GLY:N	2.41	0.40
1:B:511:PHE:HE1	1:B:551:PRO:HD2	1.85	0.40
1:A:209:LYS:C	1:A:211:PHE:H	2.29	0.40
1:B:146:ARG:NH2	1:B:164:GLU:OE1	2.54	0.40
1:C:408:ARG:HD3	1:C:409:LEU:HB2	2.04	0.40
2:E:6:GLU:HB2	2:E:90:ARG:HH21	1.86	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	609/727 (84%)	536 (88%)	69 (11%)	4 (1%)	18 51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	619/727 (85%)	544 (88%)	73 (12%)	2 (0%)	36	65
1	C	640/727 (88%)	557 (87%)	79 (12%)	4 (1%)	21	54
1	D	611/727 (84%)	522 (85%)	88 (14%)	1 (0%)	43	72
2	E	146/149 (98%)	130 (89%)	16 (11%)	0	100	100
All	All	2625/3057 (86%)	2289 (87%)	325 (12%)	11 (0%)	31	61

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	210	THR
1	B	638	LEU
1	C	209	LYS
1	A	288	ASP
1	C	288	ASP
1	D	210	THR
1	A	210	THR
1	C	379	PRO
1	B	379	PRO
1	A	379	PRO
1	A	207	PRO

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	528/630 (84%)	519 (98%)	9 (2%)	53	69
1	B	537/630 (85%)	533 (99%)	4 (1%)	76	78
1	C	557/630 (88%)	550 (99%)	7 (1%)	61	72
1	D	530/630 (84%)	526 (99%)	4 (1%)	73	77
2	E	126/127 (99%)	126 (100%)	0	100	100
All	All	2278/2647 (86%)	2254 (99%)	24 (1%)	63	74

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	LEU
1	A	199	VAL
1	A	228	LEU
1	A	451	VAL
1	A	495	LEU
1	A	519	LEU
1	A	572	LEU
1	A	611	LEU
1	A	612	TRP
1	B	451	VAL
1	B	495	LEU
1	B	519	LEU
1	B	572	LEU
1	C	210	THR
1	C	451	VAL
1	C	479	ILE
1	C	519	LEU
1	C	568	LEU
1	C	572	LEU
1	C	643	ILE
1	D	39	LEU
1	D	451	VAL
1	D	519	LEU
1	D	611	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	118	GLN
1	A	127	ASN
1	A	217	ASN
1	A	237	ASN
1	A	251	ASN
1	A	257	HIS
1	A	264	HIS
1	A	364	ASN
1	A	368	GLN
1	A	417	GLN
1	A	571	ASN
1	A	581	HIS
1	A	586	HIS
1	B	40	GLN

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Mol	Chain	Res	Type
1	B	56	ASN
1	B	118	GLN
1	B	217	ASN
1	B	237	ASN
1	B	251	ASN
1	B	256	GLN
1	B	364	ASN
1	B	368	GLN
1	B	482	GLN
1	B	586	HIS
1	C	37	ASN
1	C	59	GLN
1	C	118	GLN
1	C	127	ASN
1	C	227	HIS
1	C	237	ASN
1	C	251	ASN
1	C	256	GLN
1	C	309	GLN
1	C	571	ASN
1	C	581	HIS
1	D	37	ASN
1	D	84	HIS
1	D	127	ASN
1	D	217	ASN
1	D	237	ASN
1	D	251	ASN
1	D	260	GLN
1	D	309	GLN
1	D	417	GLN
1	D	445	ASN
1	D	571	ASN
1	D	586	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [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.

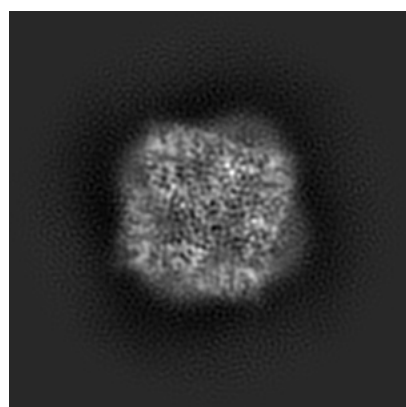
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8962. These allow visual inspection of the internal detail of the map and identification of artifacts.

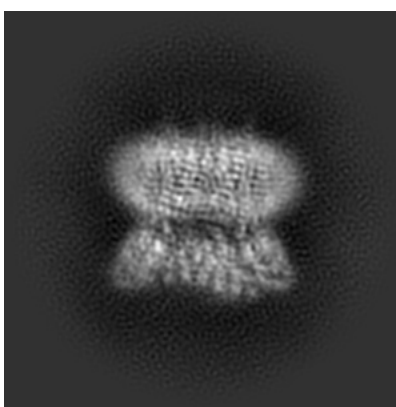
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

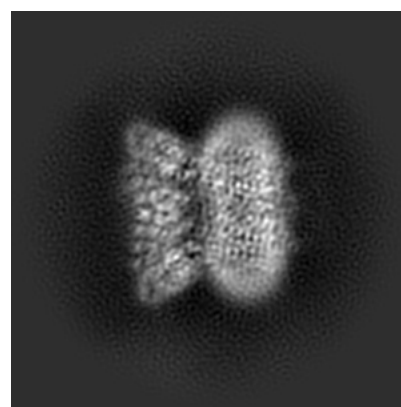
6.1.1 Primary map



X



Y

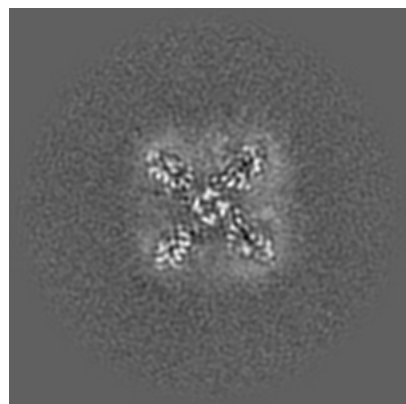


Z

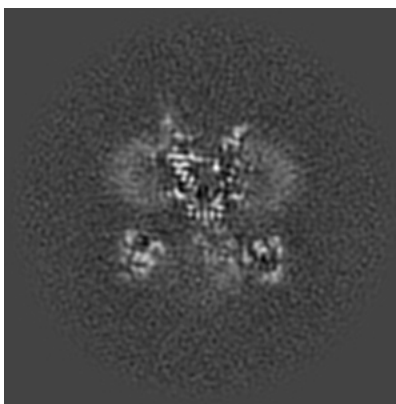
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

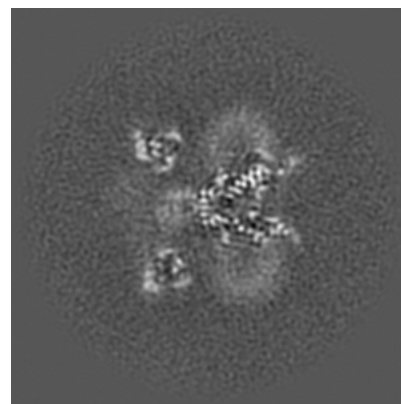
6.2.1 Primary map



X Index: 120



Y Index: 120

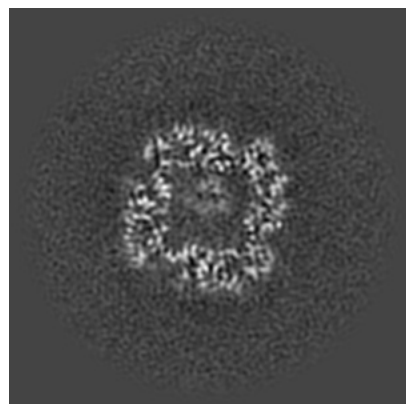


Z Index: 120

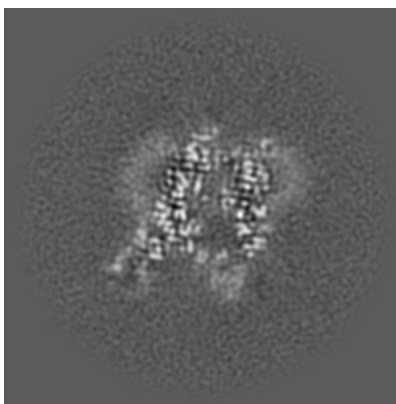
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

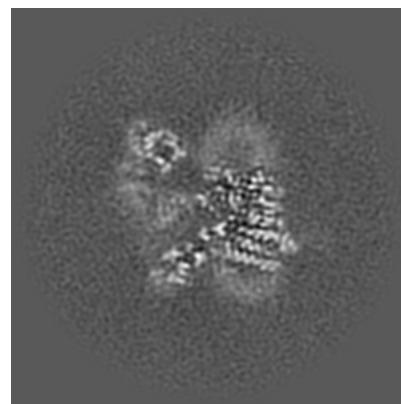
6.3.1 Primary map



X Index: 97



Y Index: 142

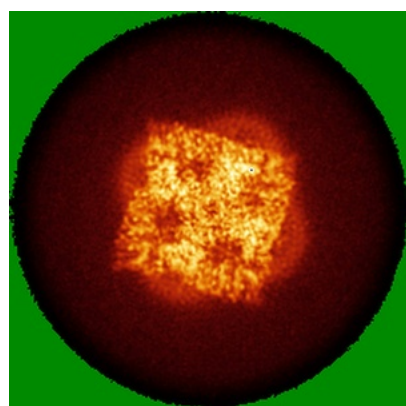


Z Index: 131

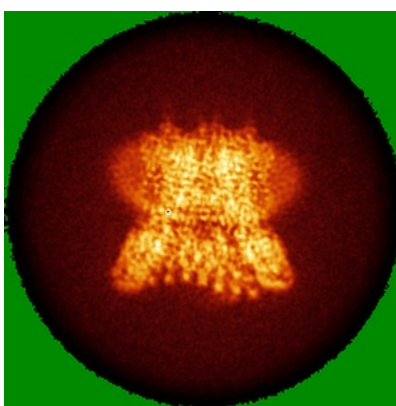
The images above show the largest variance slices of the map in three orthogonal directions.

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

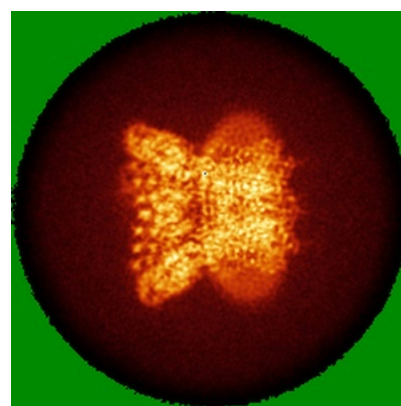
6.4.1 Primary map



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

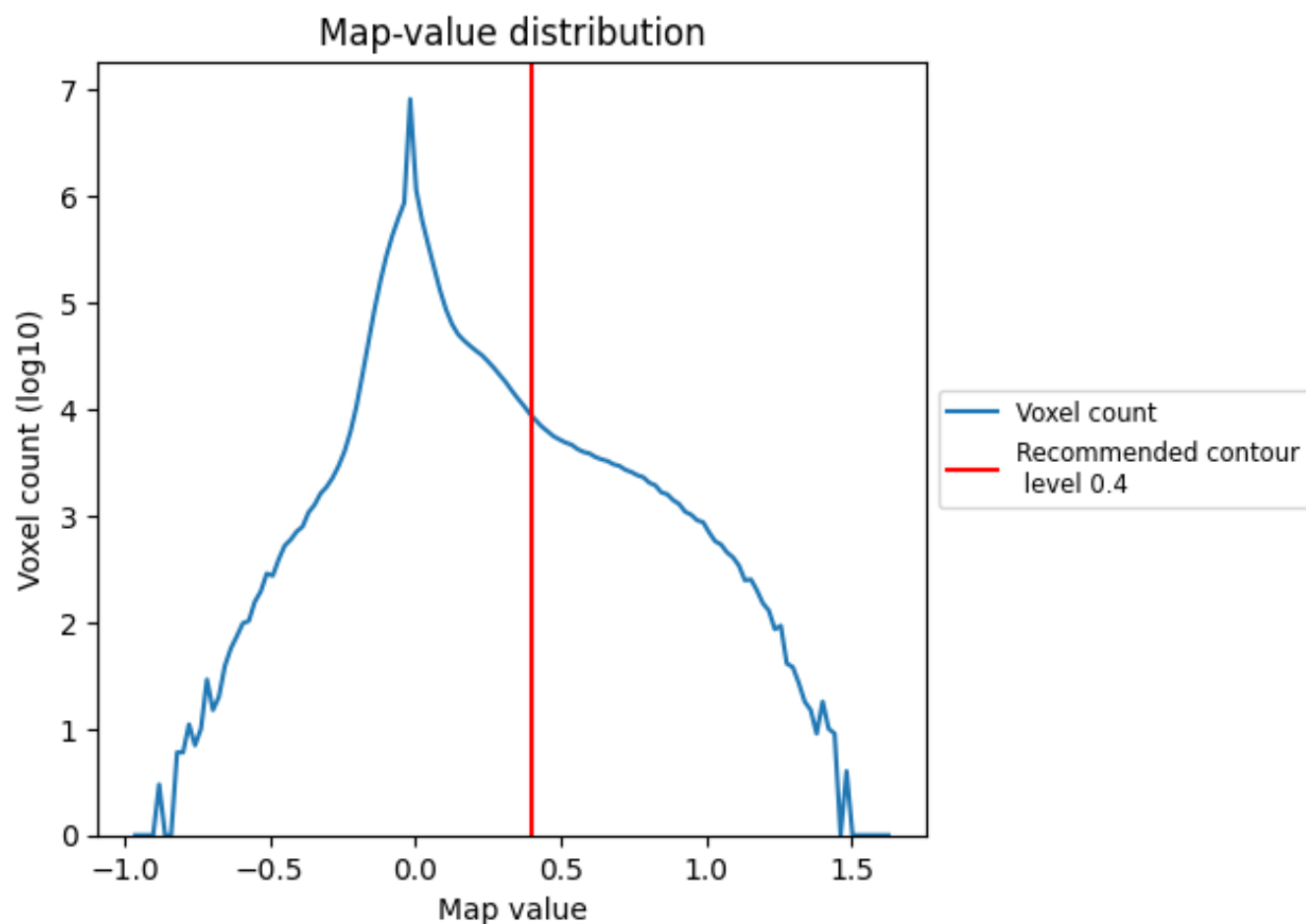
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

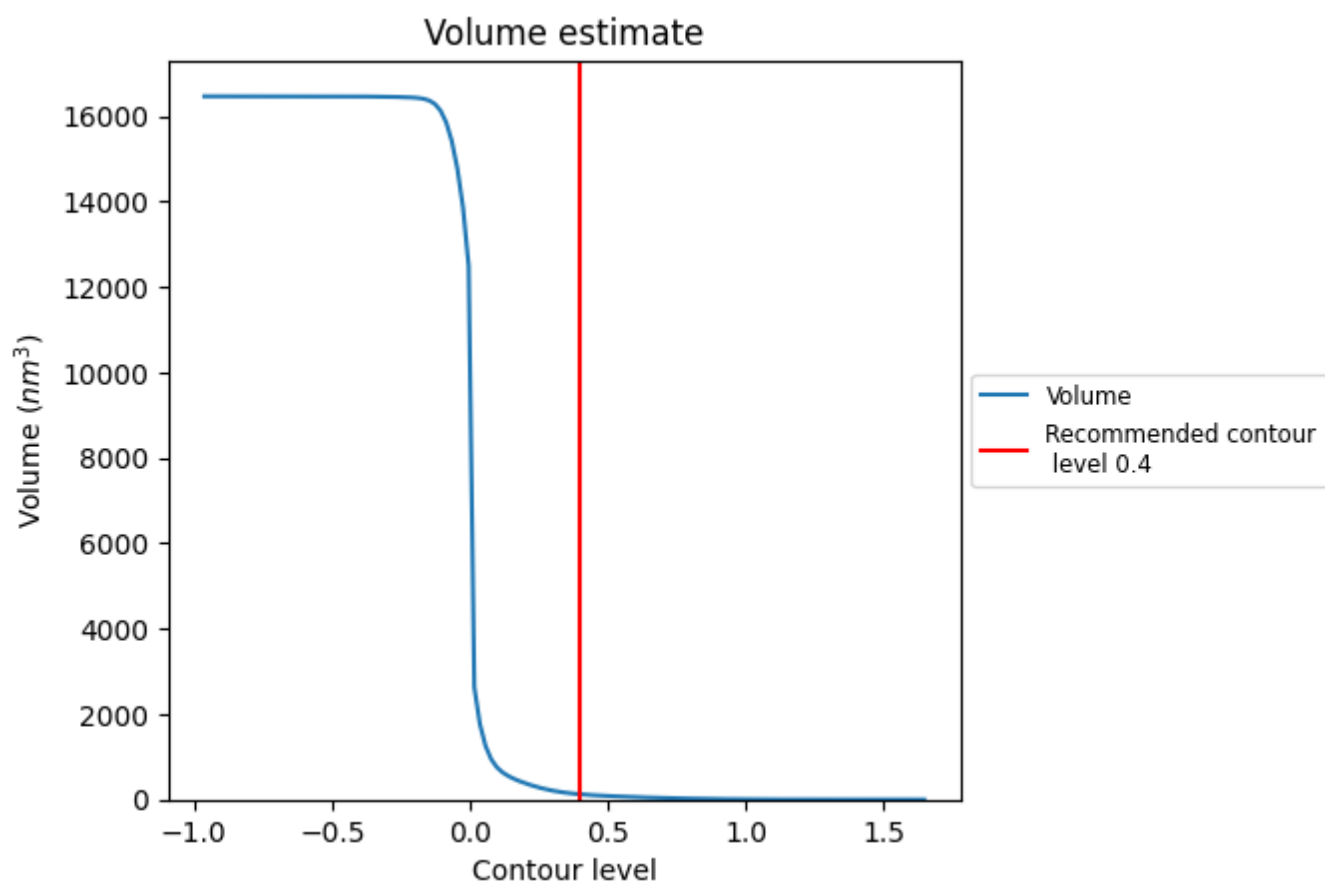
This section contains the results of statistical analysis of the map.

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.

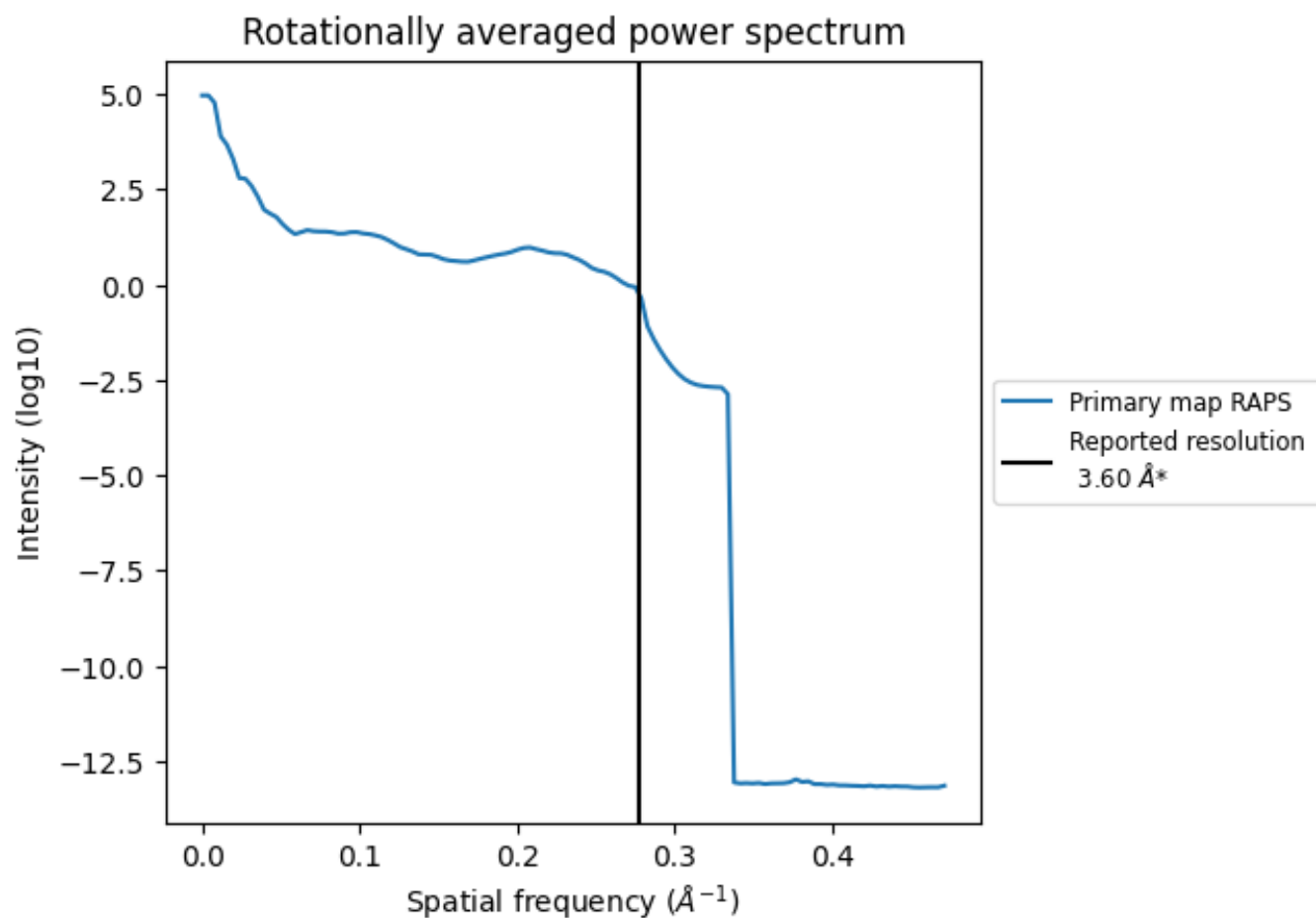
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 124 nm³; this corresponds to an approximate mass of 112 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.

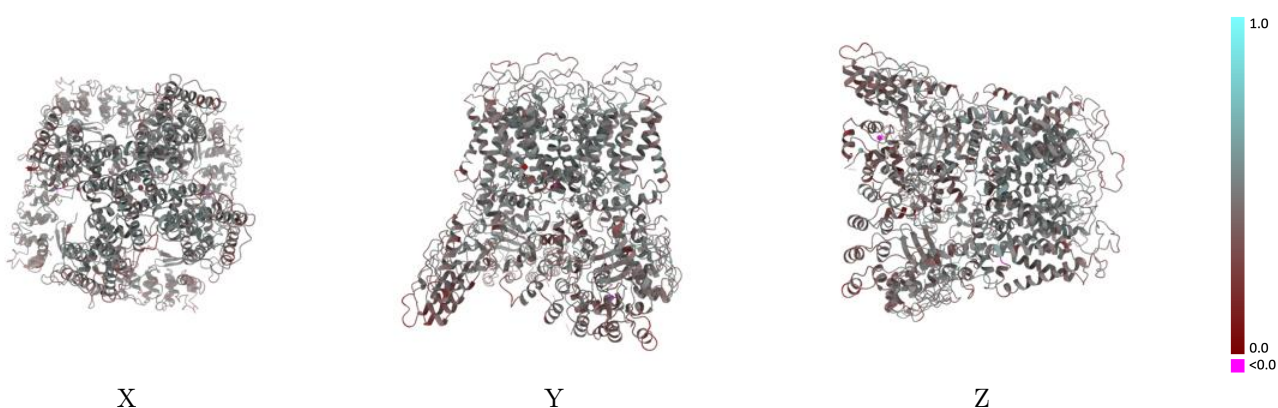
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8962 and PDB model 6E2G. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)

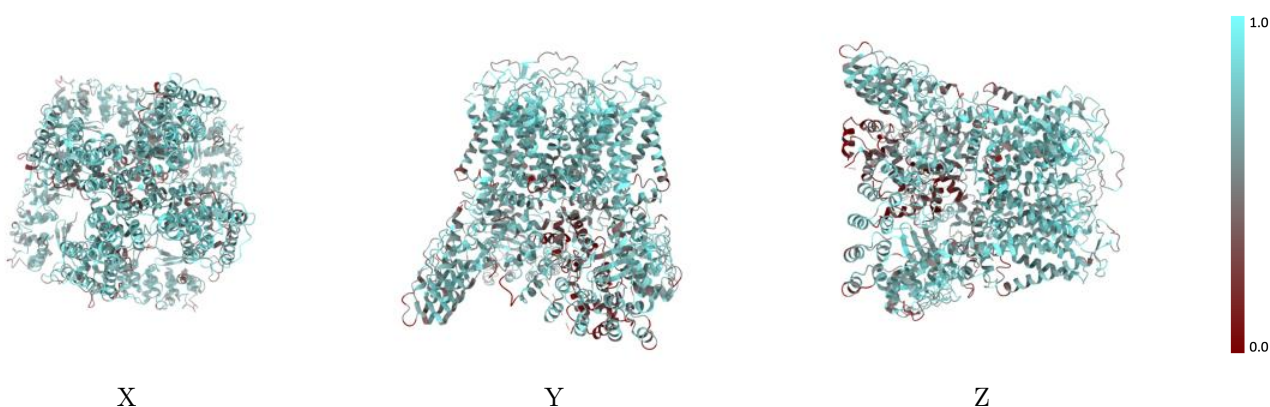
This section was not generated.

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



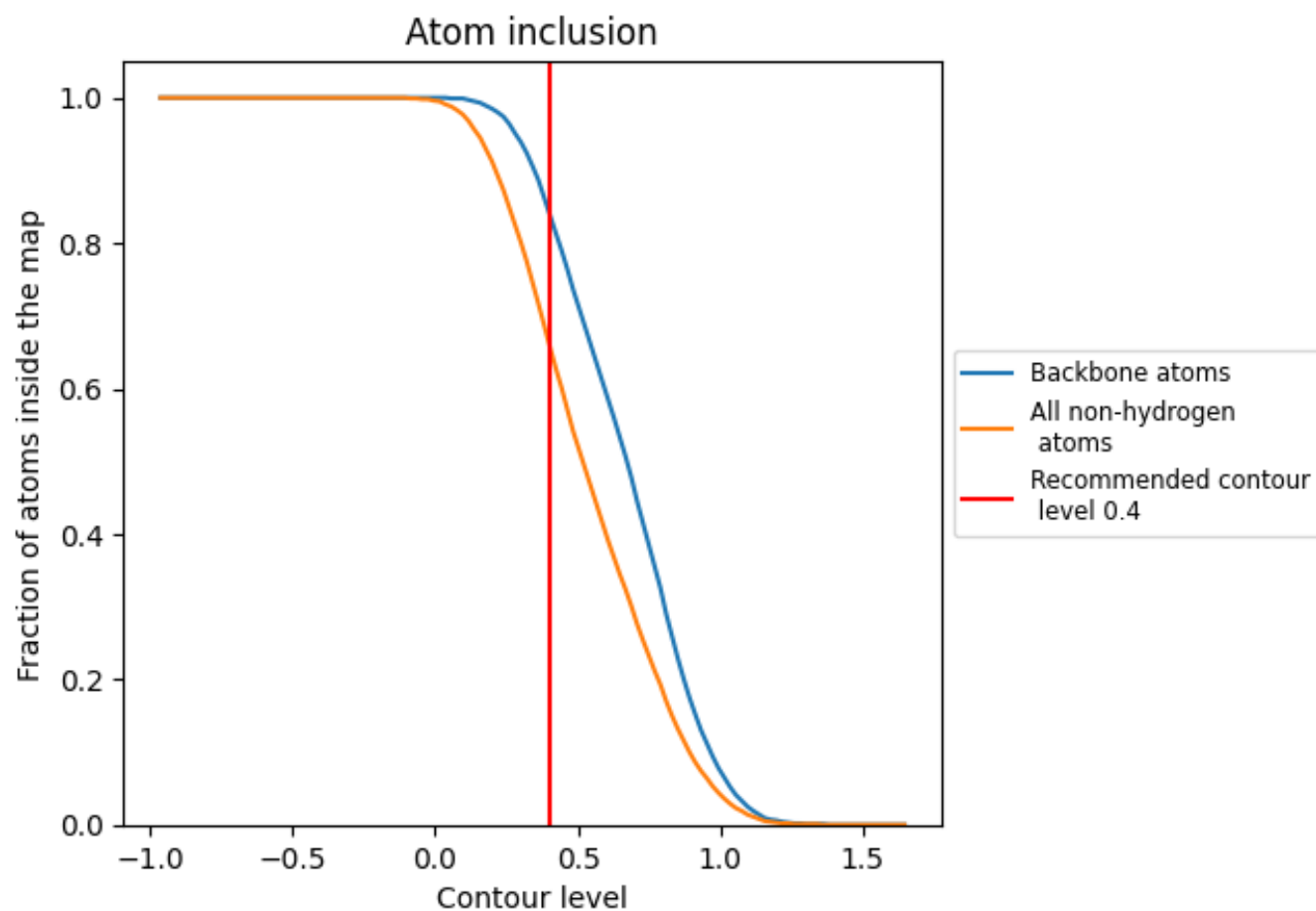
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.4).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 66% 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 (0.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6600	0.4410
A	0.6850	0.4400
B	0.6790	0.4420
C	0.6750	0.4510
D	0.6860	0.4460
E	0.3130	0.3800

