



Full wwPDB EM Validation Report ⓘ

Sep 1, 2026 – 04:34 PM EDT

PDB ID : 11AM / pdb_000011am
EMDB ID : EMD-75590
Title : Cryo-EM structure of a bovine CLC-K S68N/R355K chloride channel with 100 mM Ca²⁺
Authors : Chien, C.-T.; Chiu, W.; Maduke, M.
Deposited on : 2026-02-13
Resolution : 3.40 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.dev133
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) : 2.50

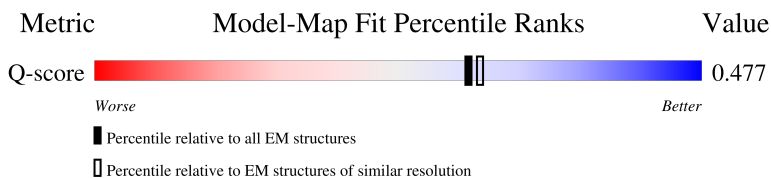
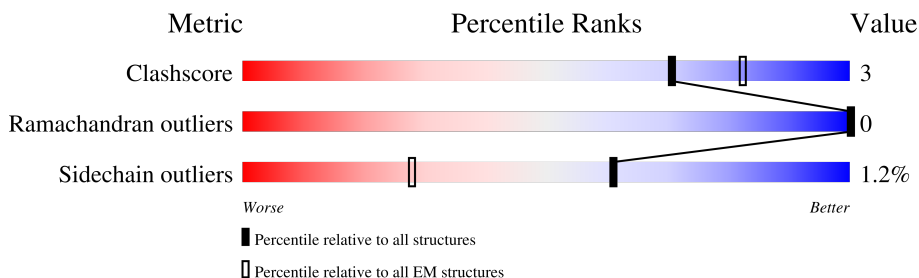
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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	14717 (2.90 - 3.90)

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	A	671	12% 76% 16% 7%
1	B	671	12% 76% 16% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19308 atoms, of which 9718 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 Chloride voltage-gated channel Ka.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	621	Total	C	H	N	O	S	0	0
			9653	3145	4859	778	846	25		
1	B	621	Total	C	H	N	O	S	0	0
			9653	3145	4859	778	846	25		

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	MET	-	initiating methionine	UNP E1B792
A	68	ASN	SER	engineered mutation	UNP E1B792
A	355	LYS	ARG	engineered mutation	UNP E1B792
A	373	GLN	ASN	engineered mutation	UNP E1B792
A	688	SER	-	expression tag	UNP E1B792
A	689	ASN	-	expression tag	UNP E1B792
A	690	SER	-	expression tag	UNP E1B792
A	691	LEU	-	expression tag	UNP E1B792
A	692	GLU	-	expression tag	UNP E1B792
A	693	VAL	-	expression tag	UNP E1B792
A	694	LEU	-	expression tag	UNP E1B792
A	695	PHE	-	expression tag	UNP E1B792
A	696	GLN	-	expression tag	UNP E1B792
B	26	MET	-	initiating methionine	UNP E1B792
B	68	ASN	SER	engineered mutation	UNP E1B792
B	355	LYS	ARG	engineered mutation	UNP E1B792
B	373	GLN	ASN	engineered mutation	UNP E1B792
B	688	SER	-	expression tag	UNP E1B792
B	689	ASN	-	expression tag	UNP E1B792
B	690	SER	-	expression tag	UNP E1B792
B	691	LEU	-	expression tag	UNP E1B792
B	692	GLU	-	expression tag	UNP E1B792
B	693	VAL	-	expression tag	UNP E1B792
B	694	LEU	-	expression tag	UNP E1B792
B	695	PHE	-	expression tag	UNP E1B792
B	696	GLN	-	expression tag	UNP E1B792

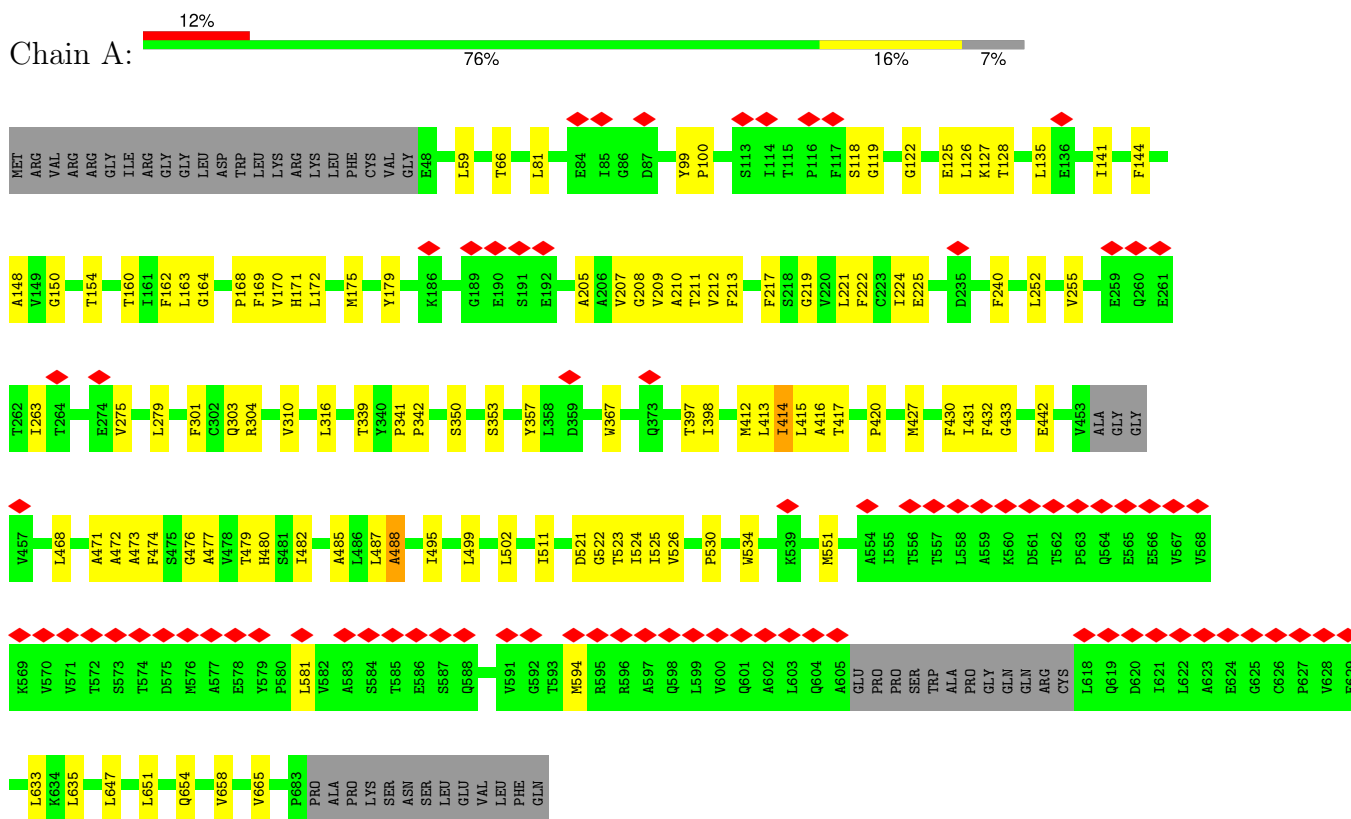
- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total 1	Cl 1	0
2	B	1	Total 1	Cl 1	0

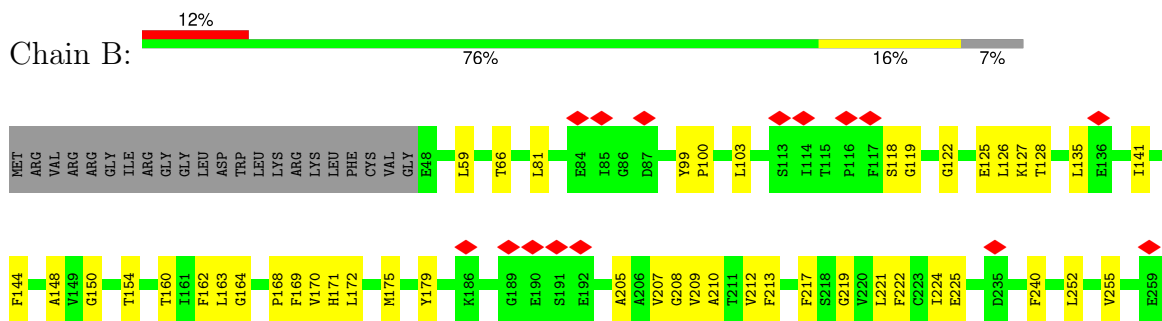
3 Residue-property plots

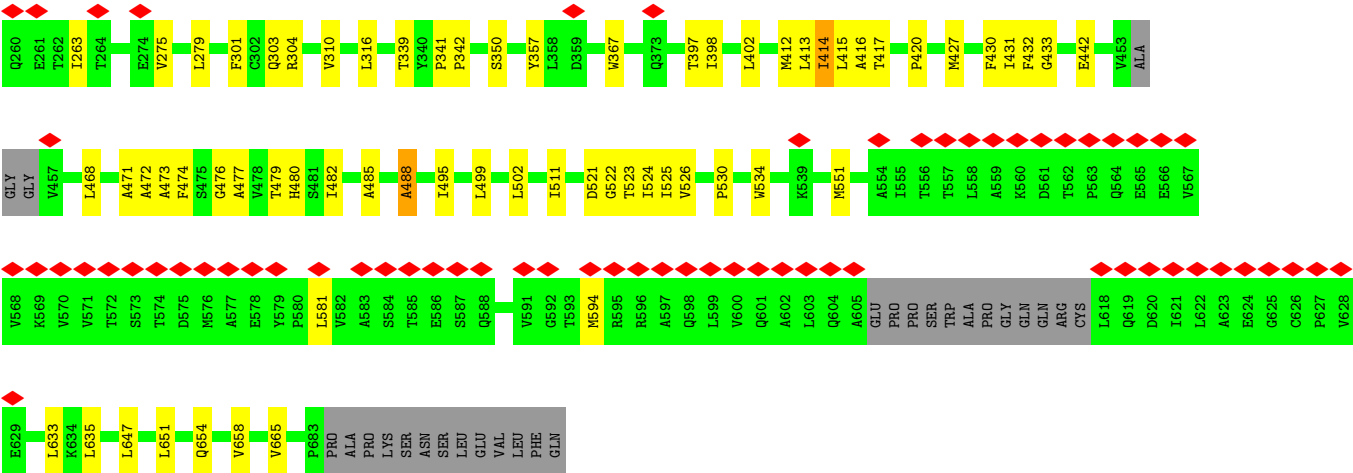
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Chloride voltage-gated channel Ka



• Molecule 1: Chloride voltage-gated channel Ka





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	191200	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.788	Depositor
Minimum map value	-0.394	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.023	Depositor
Recommended contour level	0.208	Depositor
Map size ($\text{\AA}$)	296.4, 296.4, 296.4	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.741, 0.741, 0.741	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: CL

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.16	26/4922 (0.5%)	1.23	98/6705 (1.5%)
1	B	1.17	26/4922 (0.5%)	1.24	92/6705 (1.4%)
All	All	1.16	52/9844 (0.5%)	1.23	190/13410 (1.4%)

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	171	HIS	CE1-NE2	-9.15	1.23	1.32
1	B	480	HIS	CE1-NE2	-9.02	1.23	1.32
1	B	171	HIS	CE1-NE2	-9.02	1.23	1.32
1	A	480	HIS	CE1-NE2	-9.01	1.23	1.32
1	A	480	HIS	CD2-NE2	-8.40	1.28	1.37
1	B	480	HIS	CD2-NE2	-8.30	1.28	1.37
1	A	171	HIS	CD2-NE2	-8.23	1.28	1.37
1	B	171	HIS	CD2-NE2	-8.19	1.28	1.37
1	A	304	ARG	CZ-NH2	-6.54	1.25	1.33
1	B	304	ARG	CZ-NH2	-6.54	1.25	1.33
1	A	208	GLY	N-CA	-6.32	1.38	1.45
1	B	208	GLY	N-CA	-6.31	1.38	1.45
1	B	219	GLY	N-CA	-6.23	1.38	1.45
1	A	219	GLY	N-CA	-6.23	1.38	1.45
1	A	150	GLY	N-CA	-6.20	1.38	1.45
1	B	150	GLY	N-CA	-6.20	1.38	1.45
1	B	118	SER	C-N	-6.19	1.29	1.33
1	A	118	SER	C-N	-6.19	1.29	1.33
1	A	420	PRO	CA-CB	-6.01	1.47	1.53
1	B	420	PRO	CA-CB	-5.96	1.47	1.53
1	A	433	GLY	N-CA	-5.64	1.38	1.45
1	B	433	GLY	N-CA	-5.64	1.38	1.45
1	B	476	GLY	N-CA	-5.54	1.38	1.45
1	A	476	GLY	N-CA	-5.50	1.38	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	522	GLY	N-CA	-5.46	1.38	1.45
1	B	522	GLY	N-CA	-5.46	1.38	1.45
1	A	119	GLY	CA-C	-5.40	1.48	1.52
1	B	119	GLY	CA-C	-5.39	1.48	1.52
1	B	427	MET	N-CA	-5.31	1.41	1.46
1	A	427	MET	N-CA	-5.30	1.41	1.46
1	B	477	ALA	CA-CB	-5.26	1.45	1.53
1	A	304	ARG	CD-NE	-5.25	1.38	1.46
1	B	304	ARG	CD-NE	-5.24	1.39	1.46
1	A	205	ALA	CA-CB	-5.22	1.45	1.53
1	B	205	ALA	CA-CB	-5.22	1.45	1.53
1	A	477	ALA	CA-CB	-5.21	1.45	1.53
1	B	472	ALA	CA-CB	-5.19	1.45	1.53
1	A	471	ALA	CA-CB	-5.18	1.45	1.53
1	A	119	GLY	N-CA	-5.16	1.38	1.45
1	A	485	ALA	CA-CB	-5.16	1.45	1.53
1	B	471	ALA	CA-CB	-5.16	1.45	1.53
1	A	148	ALA	CA-CB	-5.15	1.45	1.53
1	B	148	ALA	CA-CB	-5.15	1.45	1.53
1	A	416	ALA	CA-CB	-5.15	1.45	1.53
1	B	416	ALA	CA-CB	-5.15	1.45	1.53
1	A	488	ALA	CA-CB	-5.14	1.45	1.53
1	B	119	GLY	N-CA	-5.14	1.38	1.45
1	A	304	ARG	CZ-NH1	-5.10	1.25	1.32
1	B	304	ARG	CZ-NH1	-5.10	1.25	1.32
1	B	485	ALA	CA-CB	-5.10	1.45	1.53
1	B	488	ALA	CA-CB	-5.08	1.45	1.53
1	A	472	ALA	CA-CB	-5.07	1.45	1.53

All (190) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	PHE	CA-CB-CG	7.06	120.86	113.80
1	B	162	PHE	CA-CB-CG	7.06	120.86	113.80
1	A	430	PHE	CA-CB-CG	6.79	120.58	113.80
1	B	430	PHE	CA-CB-CG	6.76	120.56	113.80
1	B	213	PHE	CA-CB-CG	6.60	120.40	113.80
1	A	432	PHE	CA-CB-CG	6.55	120.35	113.80
1	B	432	PHE	CA-CB-CG	6.55	120.35	113.80
1	A	213	PHE	CA-CB-CG	6.54	120.34	113.80
1	A	480	HIS	CA-CB-CG	6.51	120.31	113.80
1	B	480	HIS	CA-CB-CG	6.50	120.30	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	PHE	CA-CB-CG	6.31	120.11	113.80
1	B	169	PHE	CA-CB-CG	6.31	120.11	113.80
1	A	525	ILE	CA-C-N	6.24	129.66	120.74
1	A	525	ILE	C-N-CA	6.24	129.66	120.74
1	A	474	PHE	CA-CB-CG	6.21	120.01	113.80
1	B	525	ILE	CA-C-N	6.20	129.60	120.74
1	B	525	ILE	C-N-CA	6.20	129.60	120.74
1	B	474	PHE	CA-CB-CG	6.13	119.93	113.80
1	B	222	PHE	CA-CB-CG	6.12	119.92	113.80
1	A	222	PHE	CA-CB-CG	6.08	119.88	113.80
1	A	521	ASP	CA-CB-CG	5.95	118.55	112.60
1	B	521	ASP	CA-CB-CG	5.95	118.55	112.60
1	B	221	LEU	CA-C-N	5.84	128.04	120.44
1	B	221	LEU	C-N-CA	5.84	128.04	120.44
1	A	534	TRP	CA-C-N	5.84	131.27	122.98
1	A	534	TRP	C-N-CA	5.84	131.27	122.98
1	B	534	TRP	CA-C-N	5.84	131.27	122.98
1	B	534	TRP	C-N-CA	5.84	131.27	122.98
1	A	221	LEU	CA-C-N	5.83	128.02	120.44
1	A	221	LEU	C-N-CA	5.83	128.02	120.44
1	A	144	PHE	CA-CB-CG	5.76	119.56	113.80
1	B	144	PHE	CA-CB-CG	5.75	119.55	113.80
1	A	171	HIS	CA-CB-CG	5.71	119.51	113.80
1	B	171	HIS	CA-CB-CG	5.67	119.47	113.80
1	A	163	LEU	CA-C-N	5.65	129.67	122.83
1	A	163	LEU	C-N-CA	5.65	129.67	122.83
1	B	163	LEU	CA-C-N	5.65	129.67	122.83
1	B	163	LEU	C-N-CA	5.65	129.67	122.83
1	A	172	LEU	CA-C-N	5.62	127.75	120.44
1	A	172	LEU	C-N-CA	5.62	127.75	120.44
1	B	172	LEU	CA-C-N	5.62	127.75	120.44
1	B	172	LEU	C-N-CA	5.62	127.75	120.44
1	A	126	LEU	CA-C-N	5.62	127.75	120.44
1	A	126	LEU	C-N-CA	5.62	127.75	120.44
1	B	126	LEU	CA-C-N	5.58	127.69	120.44
1	B	126	LEU	C-N-CA	5.58	127.69	120.44
1	A	127	LYS	CA-C-N	5.58	127.69	120.44
1	A	127	LYS	C-N-CA	5.58	127.69	120.44
1	A	217	PHE	CA-C-N	5.56	127.66	120.44
1	A	217	PHE	C-N-CA	5.56	127.66	120.44
1	B	224	ILE	CA-C-N	5.54	127.98	120.44
1	B	224	ILE	C-N-CA	5.54	127.98	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	ILE	CA-C-N	5.54	127.97	120.44
1	A	224	ILE	C-N-CA	5.54	127.97	120.44
1	B	127	LYS	CA-C-N	5.53	127.63	120.44
1	B	127	LYS	C-N-CA	5.53	127.63	120.44
1	B	217	PHE	CA-C-N	5.50	127.59	120.44
1	B	217	PHE	C-N-CA	5.50	127.59	120.44
1	A	530	PRO	CA-C-N	5.50	131.10	123.07
1	A	530	PRO	C-N-CA	5.50	131.10	123.07
1	B	530	PRO	CA-C-N	5.50	131.10	123.07
1	B	530	PRO	C-N-CA	5.50	131.10	123.07
1	A	430	PHE	CA-C-N	5.50	127.99	120.46
1	A	430	PHE	C-N-CA	5.50	127.99	120.46
1	B	168	PRO	CA-C-N	5.46	127.60	120.28
1	B	168	PRO	C-N-CA	5.46	127.60	120.28
1	B	430	PHE	CA-C-N	5.46	127.95	120.46
1	B	430	PHE	C-N-CA	5.46	127.95	120.46
1	B	103	LEU	CA-C-N	5.45	127.43	120.56
1	B	103	LEU	C-N-CA	5.45	127.43	120.56
1	B	122	GLY	CA-C-N	5.45	123.67	120.24
1	B	122	GLY	C-N-CA	5.45	123.67	120.24
1	A	301	PHE	CA-CB-CG	5.39	119.19	113.80
1	A	128	THR	CA-C-N	5.38	127.84	120.46
1	A	128	THR	C-N-CA	5.38	127.84	120.46
1	B	511	ILE	CA-C-N	5.38	127.93	120.29
1	B	511	ILE	C-N-CA	5.38	127.93	120.29
1	A	511	ILE	CA-C-N	5.38	127.92	120.29
1	A	511	ILE	C-N-CA	5.38	127.92	120.29
1	A	524	ILE	CA-C-N	5.37	127.81	120.46
1	A	524	ILE	C-N-CA	5.37	127.81	120.46
1	B	301	PHE	CA-CB-CG	5.36	119.16	113.80
1	B	128	THR	CA-C-N	5.33	127.77	120.46
1	B	128	THR	C-N-CA	5.33	127.77	120.46
1	B	524	ILE	CA-C-N	5.32	127.75	120.46
1	B	524	ILE	C-N-CA	5.32	127.75	120.46
1	A	160	THR	CA-C-N	5.31	129.94	121.34
1	A	160	THR	C-N-CA	5.31	129.94	121.34
1	B	160	THR	CA-C-N	5.29	129.92	121.34
1	B	160	THR	C-N-CA	5.29	129.92	121.34
1	A	170	VAL	N-CA-CB	5.29	116.73	110.55
1	B	170	VAL	N-CA-CB	5.29	116.73	110.55
1	A	154	THR	CA-C-N	5.27	127.34	120.28
1	A	154	THR	C-N-CA	5.27	127.34	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	154	THR	CA-C-N	5.27	127.34	120.28
1	B	154	THR	C-N-CA	5.27	127.34	120.28
1	A	122	GLY	CA-C-N	5.26	123.56	120.24
1	A	122	GLY	C-N-CA	5.26	123.56	120.24
1	B	526	VAL	CA-C-N	5.25	127.84	120.28
1	B	526	VAL	C-N-CA	5.25	127.84	120.28
1	A	431	ILE	CA-C-N	5.24	127.25	120.44
1	A	431	ILE	C-N-CA	5.24	127.25	120.44
1	B	431	ILE	CA-C-N	5.24	127.25	120.44
1	B	431	ILE	C-N-CA	5.24	127.25	120.44
1	A	66	THR	CA-C-N	5.24	127.25	120.44
1	A	66	THR	C-N-CA	5.24	127.25	120.44
1	A	526	VAL	CA-C-N	5.24	127.82	120.28
1	A	526	VAL	C-N-CA	5.24	127.82	120.28
1	B	66	THR	CA-C-N	5.24	127.25	120.44
1	B	66	THR	C-N-CA	5.24	127.25	120.44
1	B	476	GLY	CA-C-N	5.24	127.25	120.44
1	B	476	GLY	C-N-CA	5.24	127.25	120.44
1	A	303	GLN	CA-C-N	5.23	127.55	120.38
1	A	303	GLN	C-N-CA	5.23	127.55	120.38
1	A	476	GLY	CA-C-N	5.22	127.23	120.44
1	A	476	GLY	C-N-CA	5.22	127.23	120.44
1	A	59	LEU	CA-C-N	5.21	127.21	120.44
1	A	59	LEU	C-N-CA	5.21	127.21	120.44
1	A	209	VAL	N-CA-CB	5.21	116.65	110.55
1	B	303	GLN	CA-C-N	5.20	127.50	120.38
1	B	303	GLN	C-N-CA	5.20	127.50	120.38
1	B	59	LEU	CA-C-N	5.19	127.19	120.44
1	B	59	LEU	C-N-CA	5.19	127.19	120.44
1	A	168	PRO	CA-C-N	5.19	127.23	120.28
1	A	168	PRO	C-N-CA	5.19	127.23	120.28
1	A	414	ILE	CA-C-N	5.19	127.66	120.29
1	A	414	ILE	C-N-CA	5.19	127.66	120.29
1	B	414	ILE	CA-C-N	5.19	127.66	120.29
1	B	414	ILE	C-N-CA	5.19	127.66	120.29
1	A	125	GLU	CA-C-N	5.18	127.18	120.44
1	A	125	GLU	C-N-CA	5.18	127.18	120.44
1	B	209	VAL	N-CA-CB	5.18	116.61	110.55
1	A	473	ALA	CA-C-N	5.17	127.16	120.44
1	A	473	ALA	C-N-CA	5.17	127.16	120.44
1	A	523	THR	CA-C-N	5.16	127.53	120.46
1	A	523	THR	C-N-CA	5.16	127.53	120.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	523	THR	CA-C-N	5.15	127.52	120.46
1	B	523	THR	C-N-CA	5.15	127.52	120.46
1	B	471	ALA	CA-C-N	5.14	127.13	120.44
1	B	471	ALA	C-N-CA	5.14	127.13	120.44
1	A	415	LEU	CA-C-N	5.14	127.16	120.28
1	A	415	LEU	C-N-CA	5.14	127.16	120.28
1	A	170	VAL	CA-C-N	5.13	127.11	120.44
1	A	170	VAL	C-N-CA	5.13	127.11	120.44
1	B	125	GLU	CA-C-N	5.13	127.11	120.44
1	B	125	GLU	C-N-CA	5.13	127.11	120.44
1	B	170	VAL	CA-C-N	5.13	127.11	120.44
1	B	170	VAL	C-N-CA	5.13	127.11	120.44
1	B	415	LEU	CA-C-N	5.12	127.15	120.28
1	B	415	LEU	C-N-CA	5.12	127.15	120.28
1	B	473	ALA	CA-C-N	5.12	127.10	120.44
1	B	473	ALA	C-N-CA	5.12	127.10	120.44
1	A	210	ALA	CA-C-N	5.11	127.39	120.44
1	A	210	ALA	C-N-CA	5.11	127.39	120.44
1	B	210	ALA	CA-C-N	5.11	127.39	120.44
1	B	210	ALA	C-N-CA	5.11	127.39	120.44
1	A	412	MET	CA-C-N	5.11	127.39	120.44
1	A	412	MET	C-N-CA	5.11	127.39	120.44
1	A	471	ALA	CA-C-N	5.11	127.08	120.44
1	A	471	ALA	C-N-CA	5.11	127.08	120.44
1	B	412	MET	CA-C-N	5.11	127.39	120.44
1	B	412	MET	C-N-CA	5.11	127.39	120.44
1	A	164	GLY	CA-C-N	5.10	129.54	121.18
1	A	164	GLY	C-N-CA	5.10	129.54	121.18
1	A	225	GLU	CA-C-N	5.09	128.70	120.82
1	A	225	GLU	C-N-CA	5.09	128.70	120.82
1	A	482	ILE	CA-C-N	5.08	127.60	120.28
1	A	482	ILE	C-N-CA	5.08	127.60	120.28
1	A	207	VAL	CA-C-N	5.08	125.58	119.94
1	A	207	VAL	C-N-CA	5.08	125.58	119.94
1	B	207	VAL	CA-C-N	5.08	125.58	119.94
1	B	207	VAL	C-N-CA	5.08	125.58	119.94
1	A	211	THR	CA-C-N	5.07	127.05	120.56
1	A	211	THR	C-N-CA	5.07	127.05	120.56
1	A	353	SER	CA-C-N	5.06	127.06	120.28
1	A	353	SER	C-N-CA	5.06	127.06	120.28
1	A	357	TYR	CA-C-N	5.06	127.99	120.31
1	A	357	TYR	C-N-CA	5.06	127.99	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	164	GLY	CA-C-N	5.05	129.47	121.18
1	B	164	GLY	C-N-CA	5.05	129.47	121.18
1	B	225	GLU	CA-C-N	5.04	128.63	120.82
1	B	225	GLU	C-N-CA	5.04	128.63	120.82
1	A	222	PHE	CA-C-N	5.03	126.98	120.44
1	A	222	PHE	C-N-CA	5.03	126.98	120.44
1	A	487	LEU	CA-C-N	5.01	127.25	120.38
1	A	487	LEU	C-N-CA	5.01	127.25	120.38
1	B	357	TYR	CA-C-N	5.01	127.93	120.31
1	B	357	TYR	C-N-CA	5.01	127.93	120.31
1	B	482	ILE	CA-C-N	5.00	127.48	120.28
1	B	482	ILE	C-N-CA	5.00	127.48	120.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4794	4859	4856	42	0
1	B	4794	4859	4856	43	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	9590	9718	9712	67	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 3.

All (67) 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:633:LEU:HD21	1:B:647:LEU:HD11	1.68	0.75
1:A:647:LEU:HD11	1:B:633:LEU:HD21	1.68	0.74
1:B:413:LEU:O	1:B:417:THR:HG23	1.91	0.71
1:A:413:LEU:O	1:A:417:THR:HG23	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:PHE:CE1	1:B:502:LEU:HD21	2.27	0.69
1:A:240:PHE:HE1	1:B:502:LEU:HD21	1.57	0.69
1:A:502:LEU:HD21	1:B:240:PHE:CE1	2.29	0.68
1:A:279:LEU:HD11	1:B:255:VAL:HG11	1.77	0.67
1:B:468:LEU:HD11	1:B:488:ALA:HB1	1.77	0.67
1:A:255:VAL:HG11	1:B:279:LEU:HD11	1.77	0.66
1:A:468:LEU:HD11	1:A:488:ALA:HB1	1.77	0.64
1:A:263:ILE:HD12	1:B:499:LEU:HD23	1.83	0.61
1:A:339:THR:O	1:A:339:THR:HG22	2.01	0.60
1:A:499:LEU:HD23	1:B:263:ILE:HD12	1.83	0.59
1:A:502:LEU:HD21	1:B:240:PHE:HE1	1.69	0.58
1:B:339:THR:HG22	1:B:339:THR:O	2.01	0.58
1:B:581:LEU:HD13	1:B:594:MET:HE1	1.87	0.56
1:A:581:LEU:HD13	1:A:594:MET:HE1	1.87	0.56
1:A:81:LEU:HD12	1:A:81:LEU:O	2.08	0.54
1:B:495:ILE:HG22	1:B:495:ILE:O	2.08	0.54
1:B:81:LEU:HD12	1:B:81:LEU:O	2.08	0.53
1:A:397:THR:HG22	1:A:398:ILE:H	1.74	0.52
1:B:397:THR:HG22	1:B:398:ILE:H	1.74	0.52
1:A:495:ILE:HG22	1:A:495:ILE:O	2.08	0.52
1:B:581:LEU:HD13	1:B:594:MET:CE	2.42	0.50
1:A:647:LEU:HD23	1:B:651:LEU:HD11	1.93	0.50
1:A:141:ILE:HD13	1:A:179:TYR:HB2	1.94	0.50
1:A:651:LEU:HD11	1:B:647:LEU:CD2	2.41	0.50
1:A:647:LEU:CD2	1:B:651:LEU:HD11	2.41	0.49
1:A:581:LEU:HD13	1:A:594:MET:CE	2.42	0.49
1:A:310:VAL:HG22	1:A:316:LEU:HD23	1.95	0.49
1:A:651:LEU:HD11	1:B:647:LEU:HD23	1.93	0.49
1:B:141:ILE:HD13	1:B:179:TYR:HB2	1.94	0.49
1:B:310:VAL:HG22	1:B:316:LEU:HD23	1.95	0.49
1:A:502:LEU:HD11	1:B:240:PHE:CE1	2.49	0.47
1:A:479:THR:O	1:A:479:THR:HG22	2.14	0.47
1:B:479:THR:HG22	1:B:479:THR:O	2.14	0.47
1:B:495:ILE:O	1:B:495:ILE:CG2	2.62	0.47
1:A:495:ILE:O	1:A:495:ILE:CG2	2.62	0.47
1:A:339:THR:HG22	1:A:350:SER:OG	2.15	0.46
1:B:339:THR:HG22	1:B:350:SER:OG	2.15	0.46
1:B:275:VAL:HG12	1:B:275:VAL:O	2.16	0.46
1:A:275:VAL:HG12	1:A:275:VAL:O	2.16	0.46
1:A:397:THR:HG22	1:A:398:ILE:N	2.31	0.45
1:B:397:THR:HG22	1:B:398:ILE:N	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:PRO:HB2	1:B:342:PRO:HD3	1.99	0.44
1:A:367:TRP:HE3	1:A:398:ILE:HG21	1.83	0.44
1:A:341:PRO:HB2	1:A:342:PRO:HD3	1.99	0.43
1:B:99:TYR:HB3	1:B:100:PRO:HD3	1.99	0.43
1:A:99:TYR:HB3	1:A:100:PRO:HD3	1.99	0.43
1:B:310:VAL:HG22	1:B:316:LEU:CD2	2.49	0.43
1:B:367:TRP:HE3	1:B:398:ILE:HG21	1.83	0.43
1:B:397:THR:HG22	1:B:398:ILE:HD12	2.01	0.43
1:A:255:VAL:CG1	1:B:279:LEU:HD11	2.47	0.42
1:A:635:LEU:HB2	1:A:658:VAL:HG22	2.01	0.42
1:A:310:VAL:HG22	1:A:316:LEU:CD2	2.49	0.42
1:A:397:THR:HG22	1:A:398:ILE:HD12	2.01	0.42
1:A:252:LEU:C	1:A:252:LEU:HD23	2.45	0.41
1:A:654:GLN:OE1	1:A:654:GLN:N	2.53	0.41
1:B:398:ILE:HG22	1:B:402:LEU:HD12	2.02	0.41
1:A:647:LEU:CD1	1:B:633:LEU:HD21	2.46	0.41
1:B:252:LEU:HD23	1:B:252:LEU:C	2.46	0.41
1:B:635:LEU:HB2	1:B:658:VAL:HG22	2.01	0.41
1:B:654:GLN:N	1:B:654:GLN:OE1	2.53	0.41
1:A:633:LEU:HD21	1:B:647:LEU:CD1	2.46	0.40
1:B:551:MET:CE	1:B:658:VAL:HG11	2.52	0.40
1:A:551:MET:CE	1:A:658:VAL:HG11	2.52	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	615/671 (92%)	598 (97%)	17 (3%)	0	100	100
1	B	615/671 (92%)	598 (97%)	17 (3%)	0	100	100
All	All	1230/1342 (92%)	1196 (97%)	34 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	514/554 (93%)	508 (99%)	6 (1%)	63	72
1	B	514/554 (93%)	508 (99%)	6 (1%)	63	72
All	All	1028/1108 (93%)	1016 (99%)	12 (1%)	61	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	LEU
1	A	175	MET
1	A	212	VAL
1	A	414	ILE
1	A	442	GLU
1	A	665	VAL
1	B	135	LEU
1	B	175	MET
1	B	212	VAL
1	B	414	ILE
1	B	442	GLU
1	B	665	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	230	HIS
1	A	643	GLN
1	B	643	GLN

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 2 ligands modelled in this entry, 2 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.

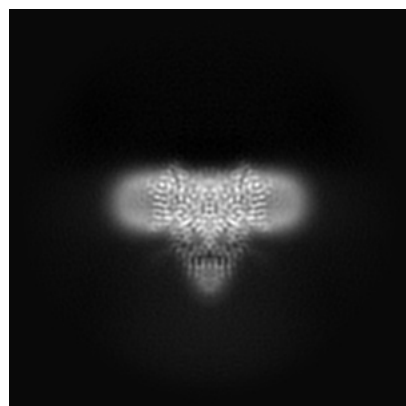
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-75590. These allow visual inspection of the internal detail of the map and identification of artifacts.

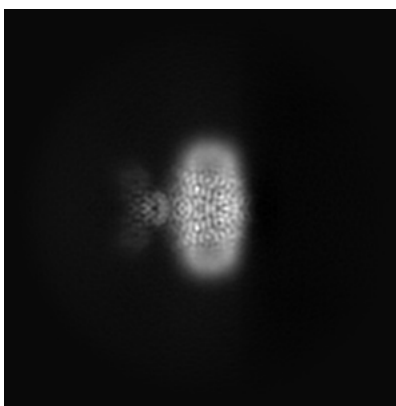
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

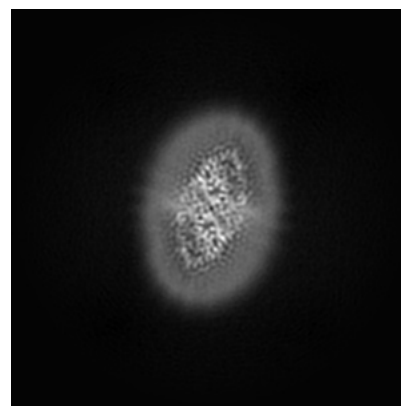
6.1.1 Primary map



X

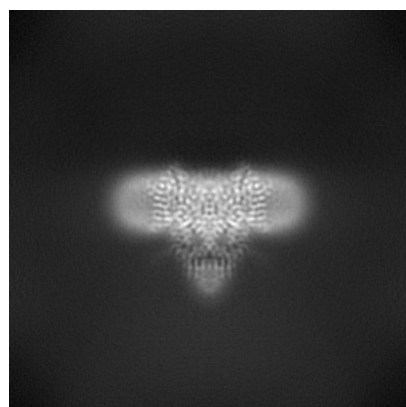


Y

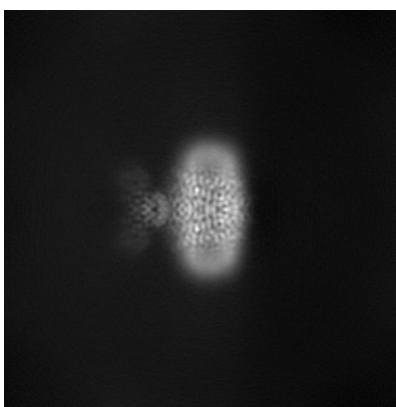


Z

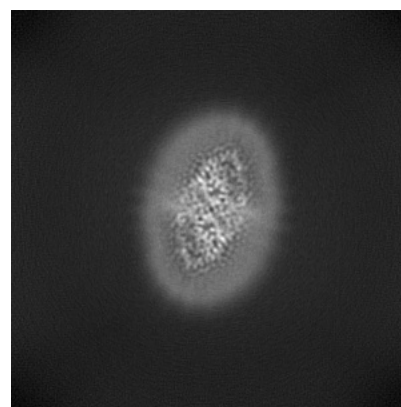
6.1.2 Raw map



X



Y

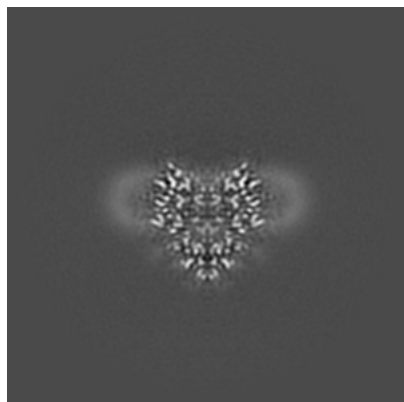


Z

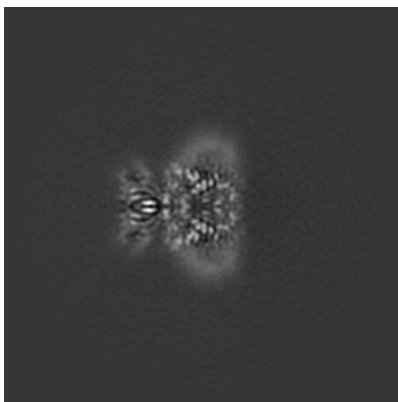
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

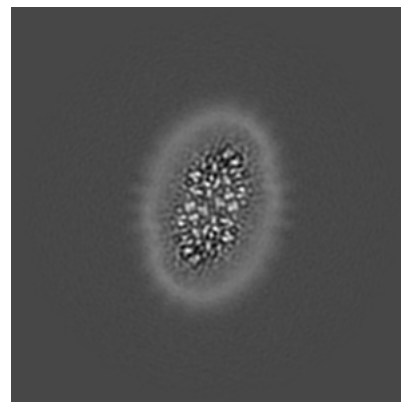
6.2.1 Primary map



X Index: 200

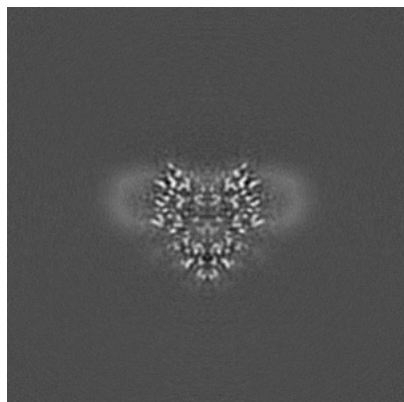


Y Index: 200

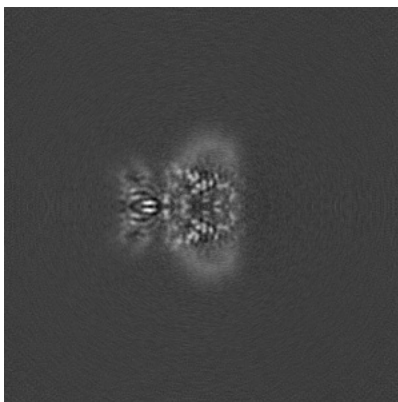


Z Index: 200

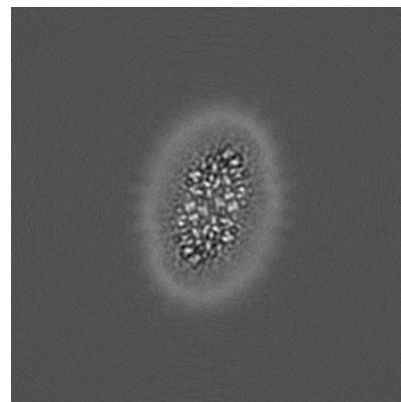
6.2.2 Raw map



X Index: 200



Y Index: 200

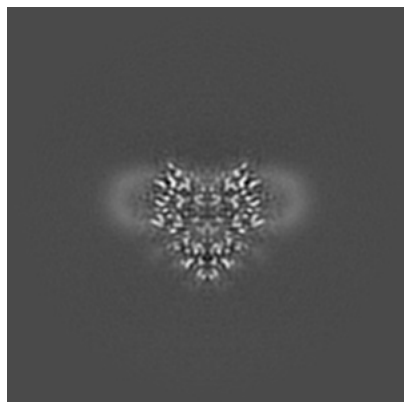


Z Index: 200

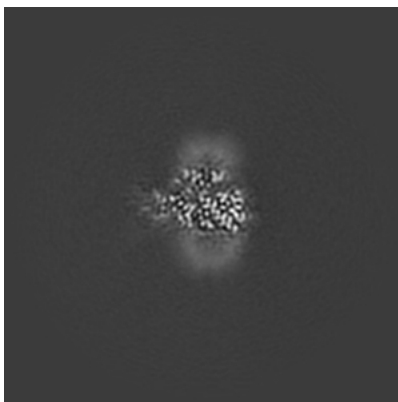
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

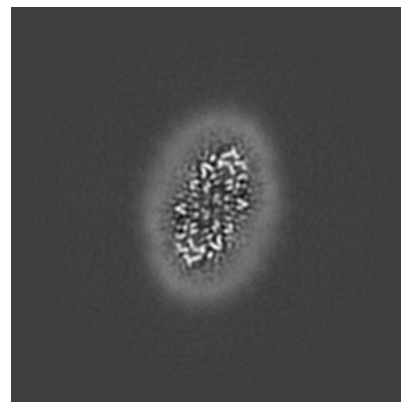
6.3.1 Primary map



X Index: 200

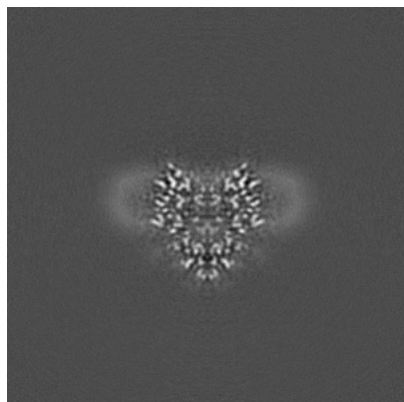


Y Index: 228

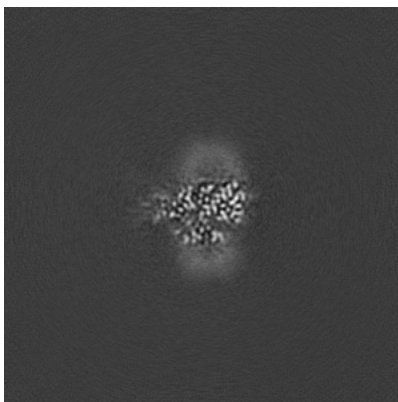


Z Index: 191

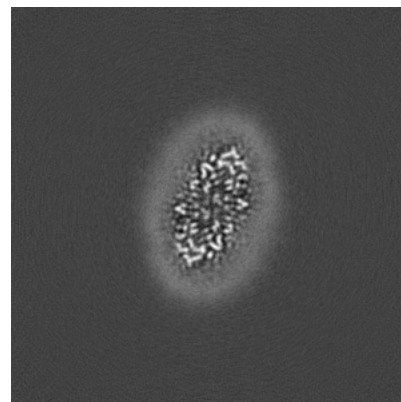
6.3.2 Raw map



X Index: 200



Y Index: 172



Z Index: 191

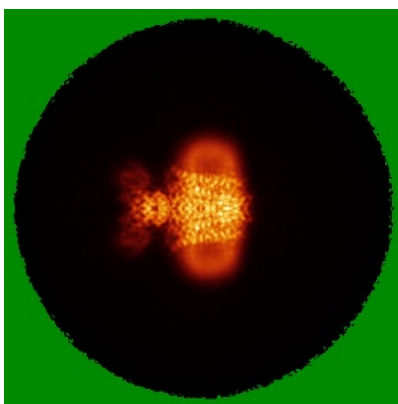
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

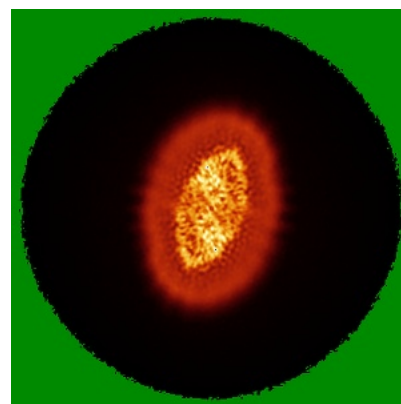
6.4.1 Primary map



X

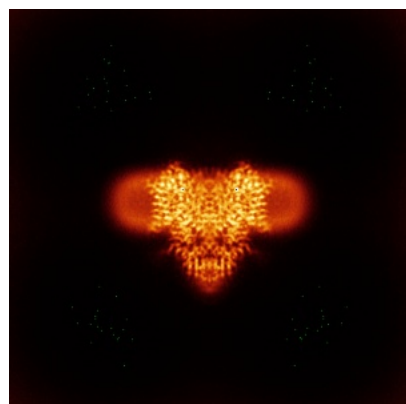


Y

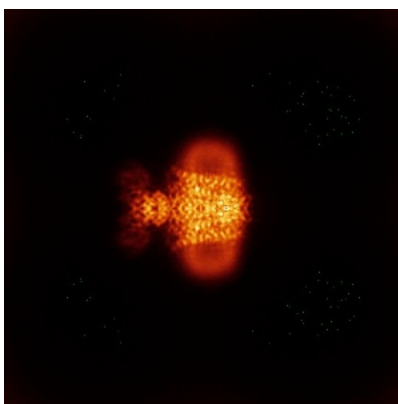


Z

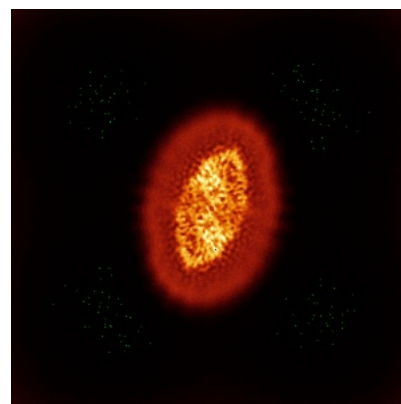
6.4.2 Raw map



X



Y

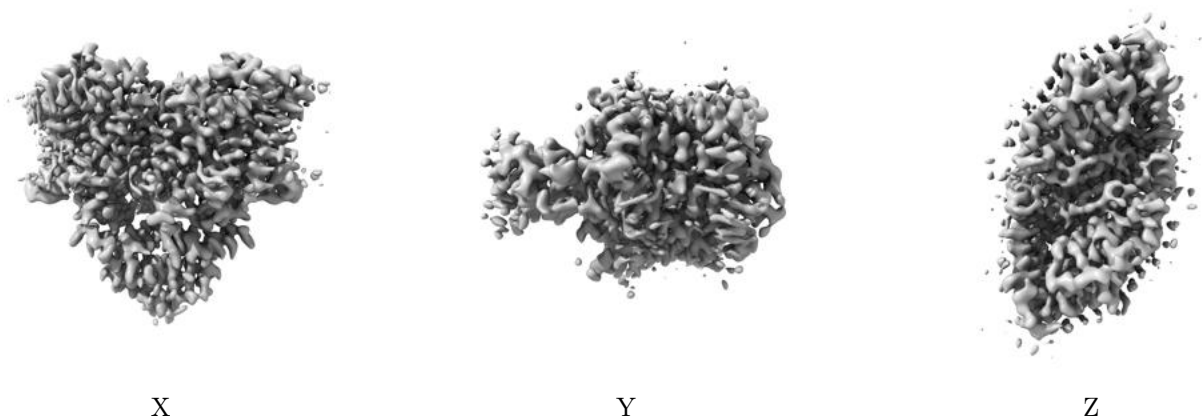


Z

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.

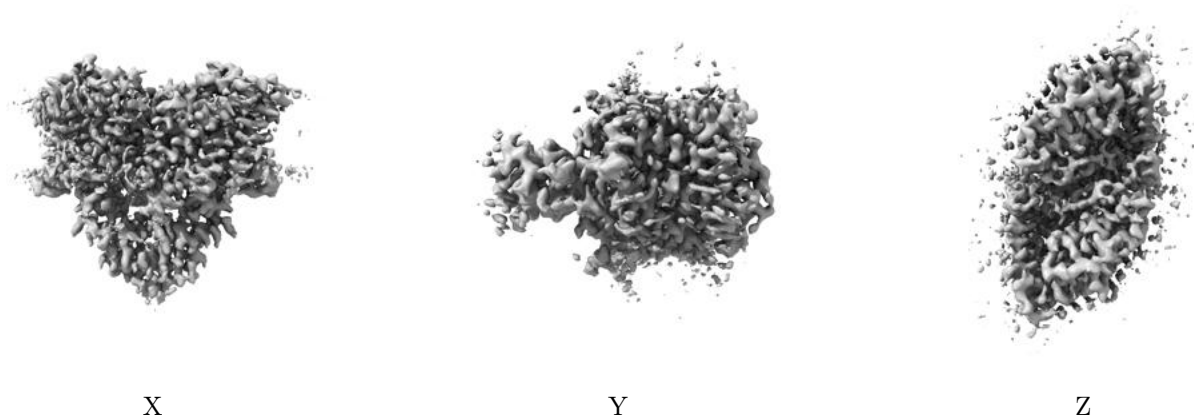
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.208. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

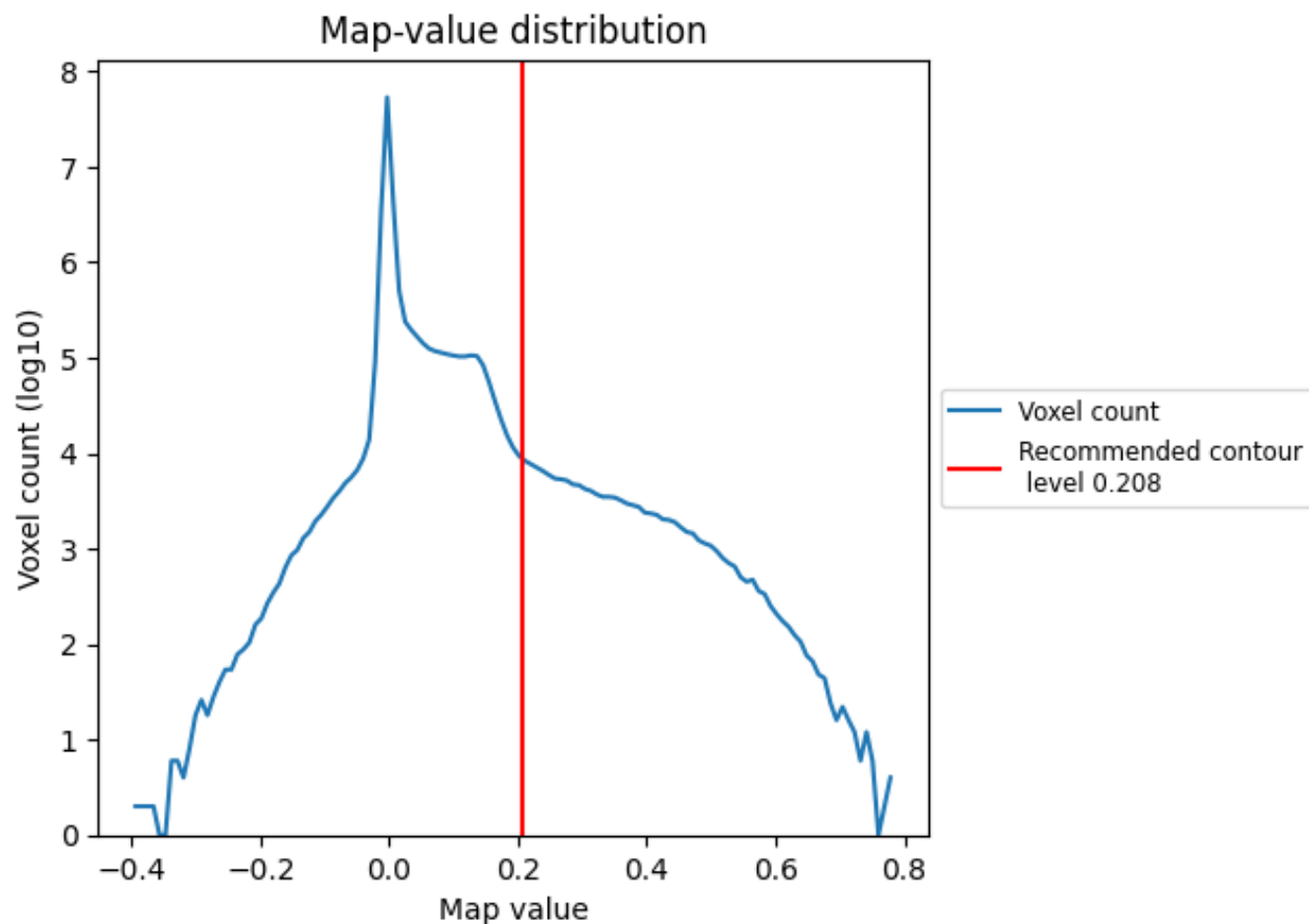
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

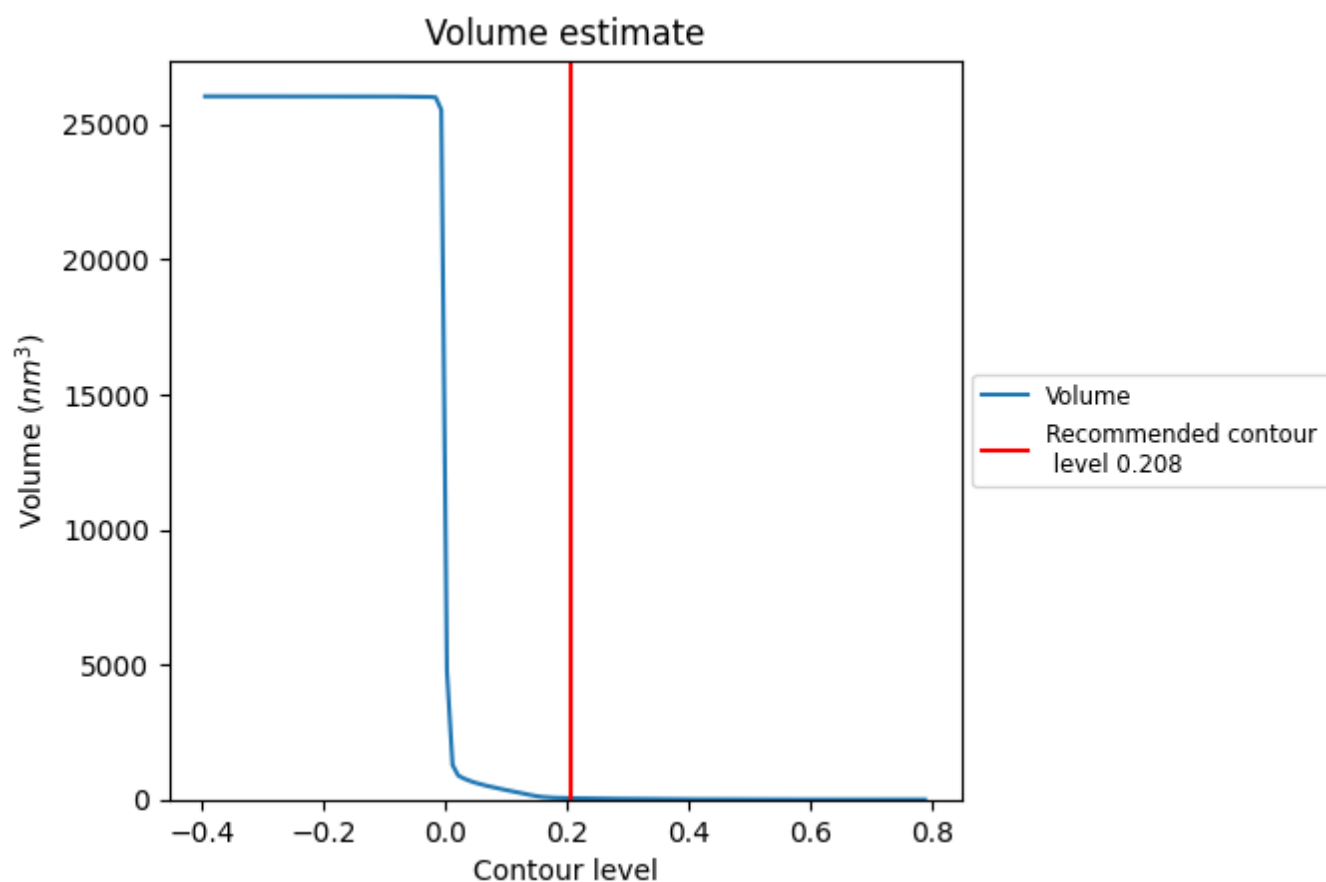
This section contains the results of statistical analysis of the map.

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.

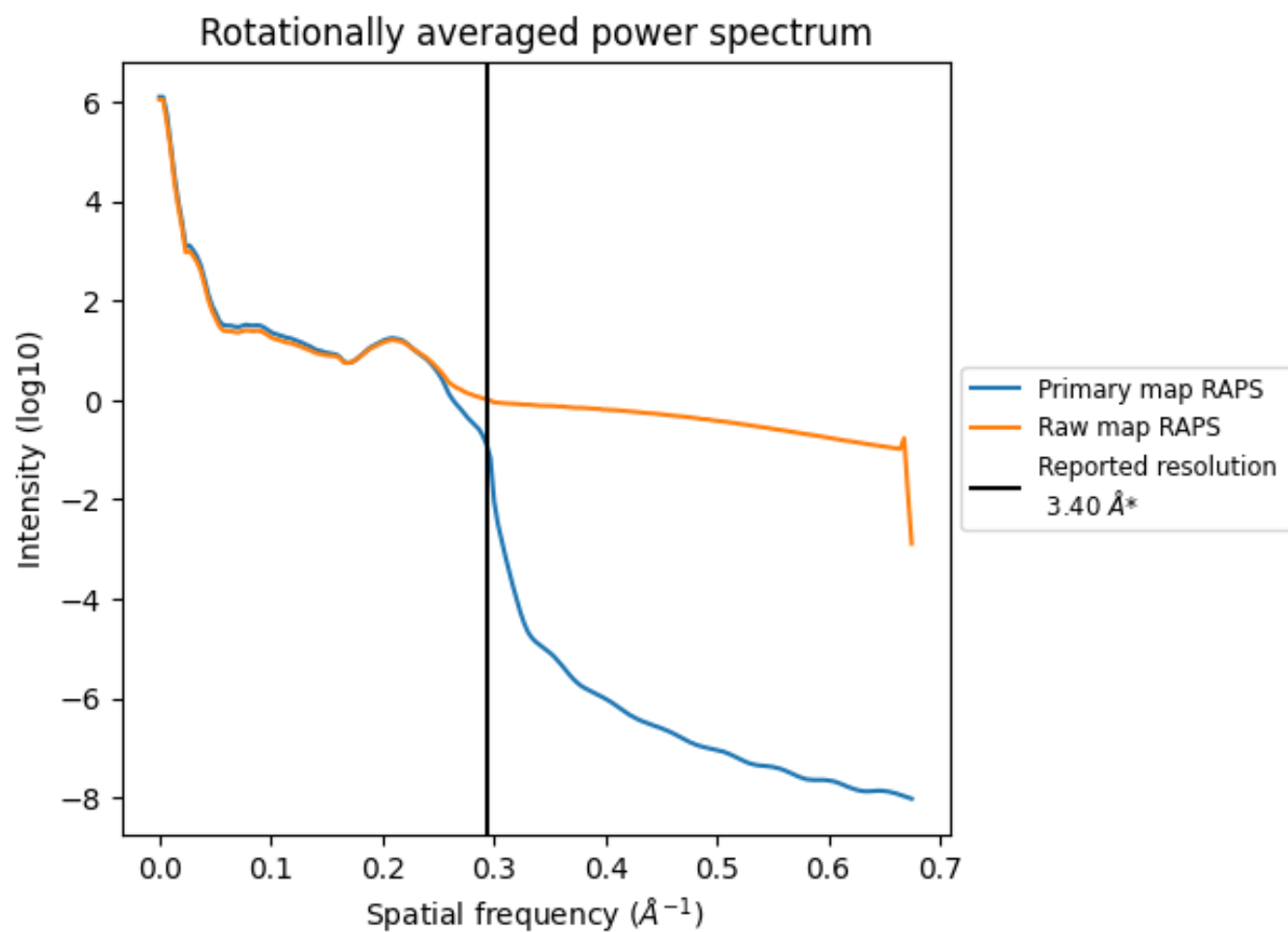
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 51 nm^3 ; this corresponds to an approximate mass of 46 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 ⓘ

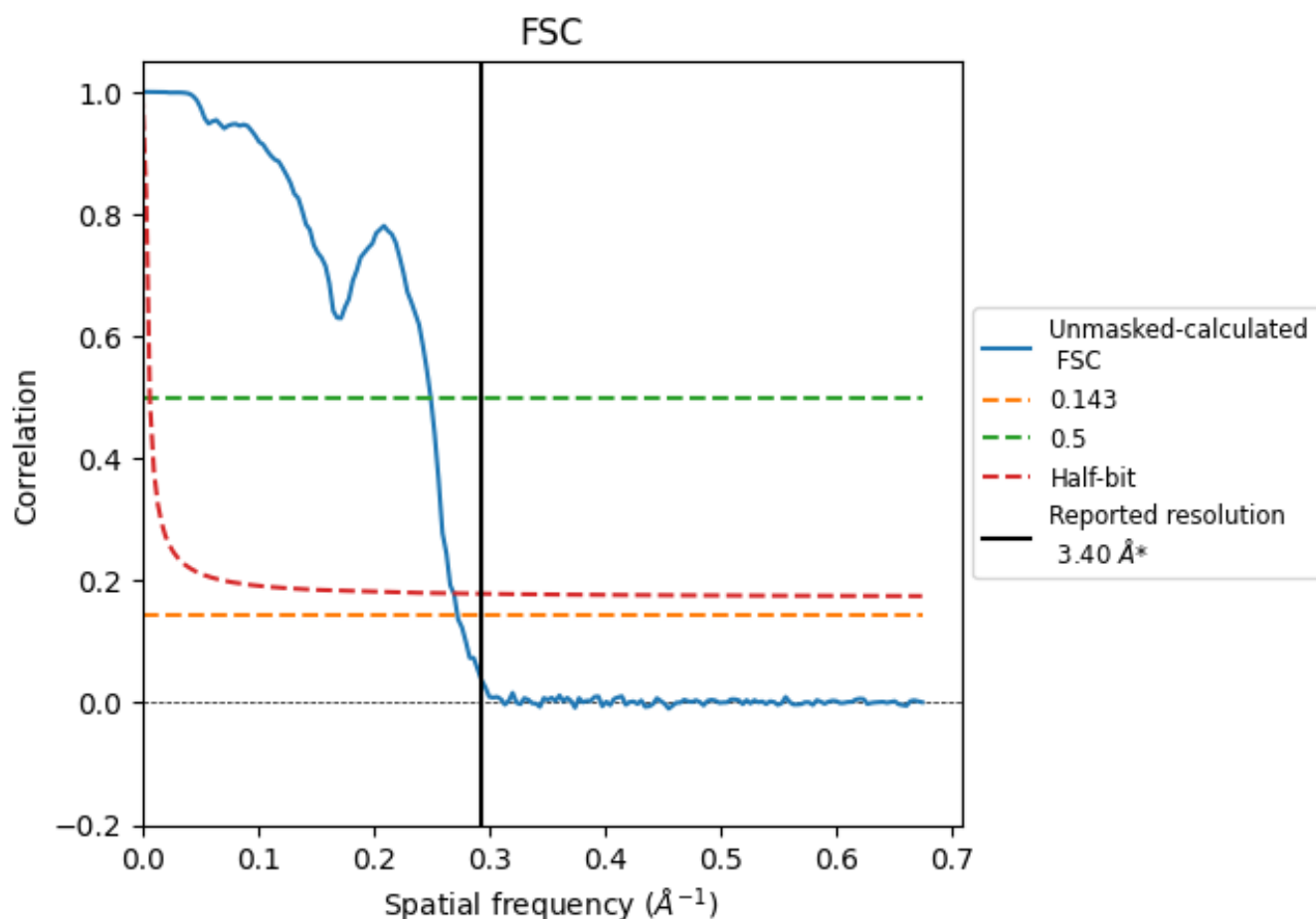


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8.2 Resolution estimates [i](#)

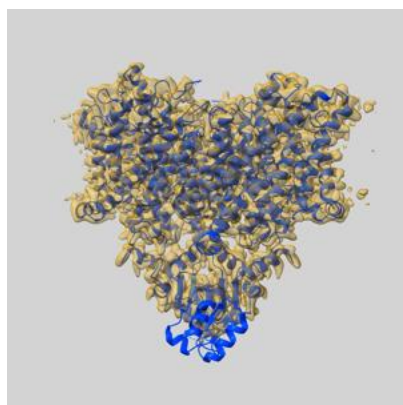
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.67	4.01	3.71

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

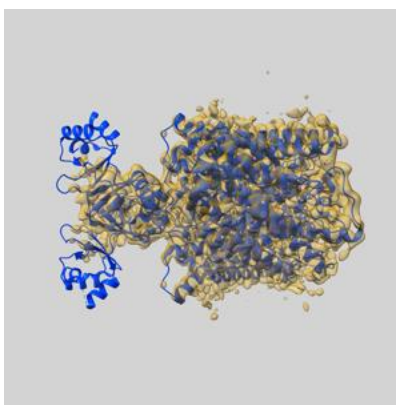
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-75590 and PDB model 11AM. Per-residue inclusion information can be found in section 3 on page 5.

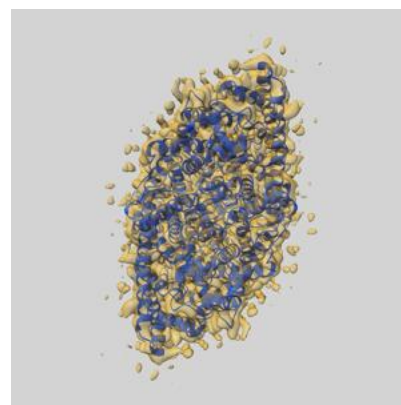
9.1 Map-model overlay [i](#)



X



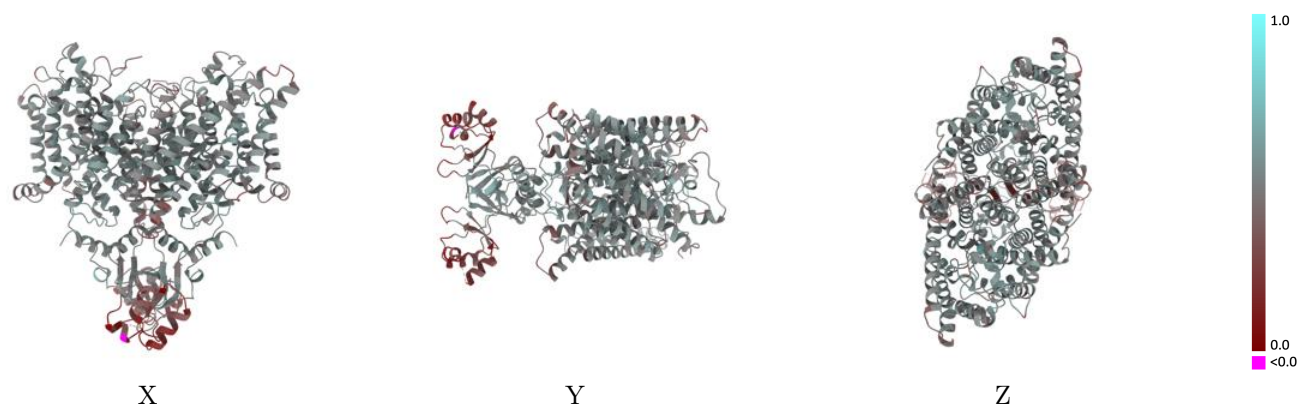
Y



Z

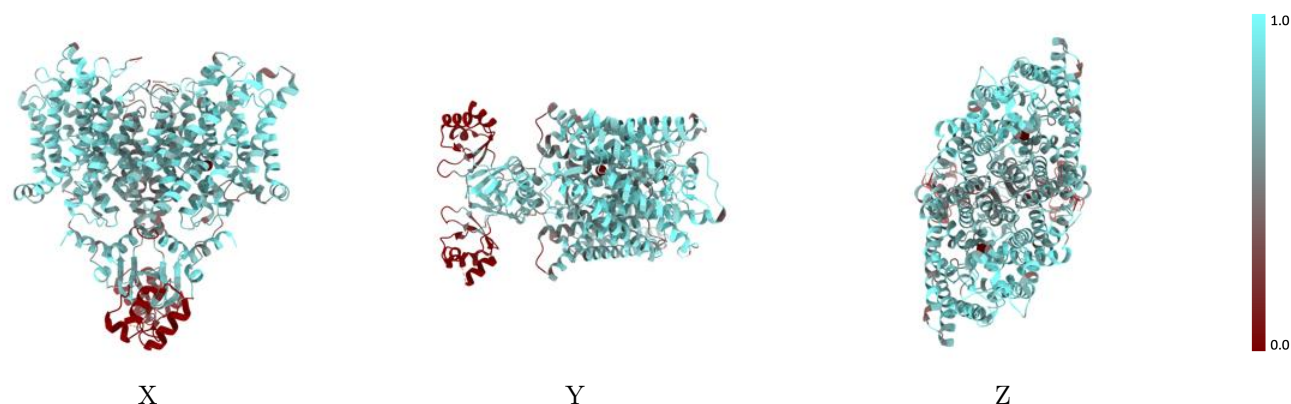
The images above show the 3D surface view of the map at the recommended contour level 0.208 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



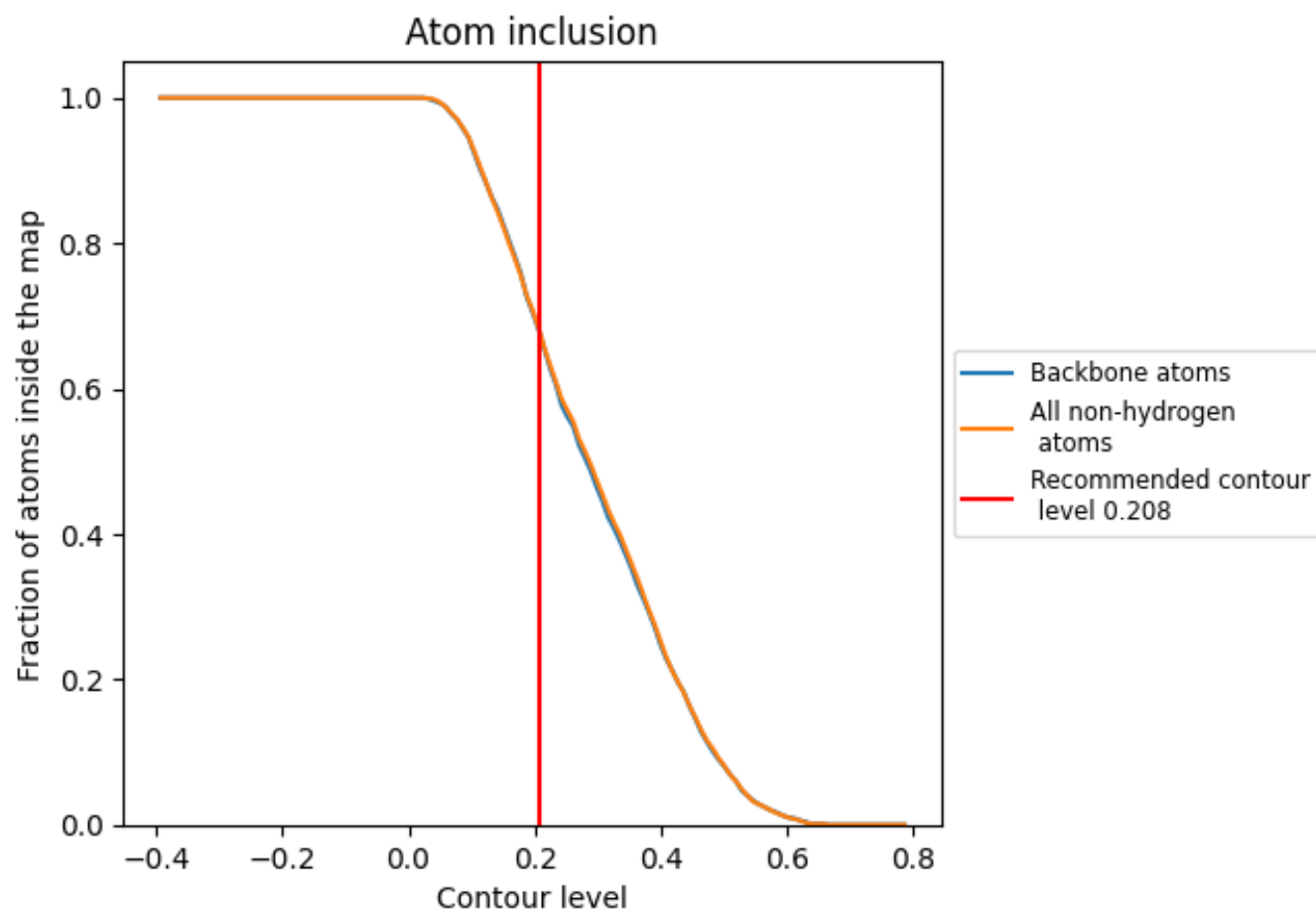
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.208).

9.4 Atom inclusion [i](#)



At the recommended contour level, 68% of all backbone atoms, 68% 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 (0.208) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6800	0.4770
A	0.6870	0.4770
B	0.6870	0.4770

