



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 02:20 AM UTC

PDB ID : 6PZT / pdb_00006pzt
EMDB ID : EMD-20537
Title : cryo-EM structure of human NKCC1
Authors : Cao, E.; Wang, Q.; Yang, X.
Deposited on : 2019-08-01
Resolution : 3.46 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

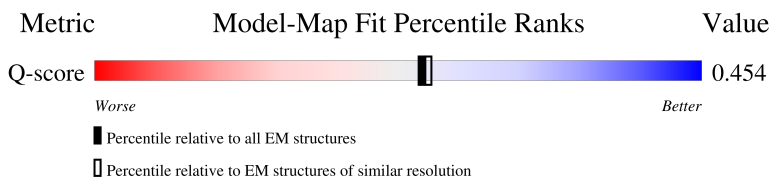
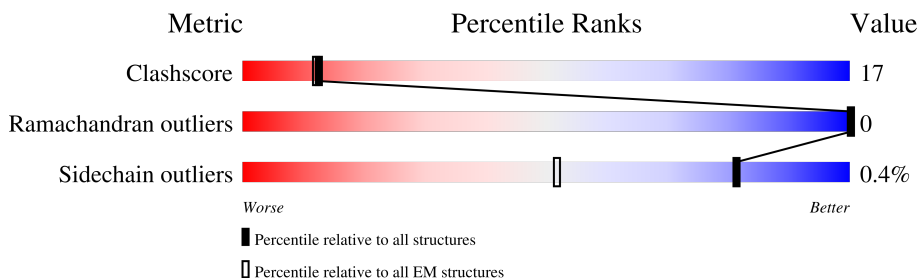
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.46 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	13788 (2.96 - 3.96)

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	1212	 27% 12% 62%
1	B	1212	 27% 12% 62%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6062 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 Solute carrier family 12 member 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	466	Total	C	N	O	S	0	0
			3031	1957	513	549	12		
1	A	466	Total	C	N	O	S	0	0
			3031	1957	513	549	12		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	289	ASN	LYS	engineered mutation	UNP P55011
B	351	ARG	GLY	engineered mutation	UNP P55011
A	289	ASN	LYS	engineered mutation	UNP P55011
A	351	ARG	GLY	engineered mutation	UNP P55011

LEU	PRO	PRO	TYR	ARG	ASP	LYS	LEU	LEU	HIS	LEU	LEU	ASP	LEU	ASP	LEU	ASP	LEU	ASP
VAL	ARG	VAL	ARG	ARG	CYS	LYS	LYS	LYS	GLN	LEU	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
ARG	ALA	LEU	LEU	LEU	ILE	SER	SER	SER	GLN	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
LYS	GLU	GLU	HIS	HIS	ARG	LYS	ARG	LYS	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	VAL
GLY	ASP	ASP	ASP	ASP	VAL	PRO	PRO	PRO	GLU	GLU	GLU	GLU	GLU	GLU	GLU	GLU	GLU	GLY
ALA	ASP	ALA	ASP	PHEN	PHEN	PRO	PRO	PRO	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	GLY
VAL	VAL	VAL	LYS	ILE	ILE	ILE	ILE	ILE	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	GLY
SER	SER	SER	GLU	GLY	GLY	VAL	VAL	SER	SER	SER	SER	SER	SER	SER	SER	SER	SER	MET
SER	GLN	GLN	GLN	GLY	GLY	PRO	PRO	GLN	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	CYS
ALA	ASP	ALA	ASP	LYS	ILE	LEU	LEU	ASN	GLN	GLN	GLN	GLN	GLN	GLN	PRO	PRO	PRO	GLY
LEU	LEU	LEU	ILE	ILE	ILE	ASN	ASN	VAL	VAL	VAL	VAL	VAL	VAL	VAL	ASN	ASN	ASN	HIS
TYR	TYR	ALA	ALA	ASN	ASN	VAL	VAL	ASN	LYS	LYS	LYS	LYS	LYS	LYS	THR	THR	THR	THR
MET	MET	ASP	ASP	ALA	ALA	ALA	ALA	ALA	SER	SER	VAL	VAL	VAL	VAL	LEU	LEU	LEU	HIS
ALA	ALA	ALA	LYS	ILE	ILE	LYS	LYS	LYS	GLY	GLY	GLY	GLY	GLY	GLY	VAL	VAL	VAL	GLY
TRP	TRP	MET	MET	ASP	ASP	GLN	GLN	SER	THR	THR	THR	THR	THR	THR	GLY	GLY	GLY	GLY
LEU	LYS	LYS	LYS	HIS	HIS	LYS	LYS	LYS	THR	THR	SER	SER	SER	SER	LEU	LEU	LEU	PRO
GLU	GLU	GLU	GLU	ASP	ASP	LEU	LEU	LEU	LYS	LYS	LEU	LEU	LEU	LEU	GLY	GLY	GLY	ARG
ALA	ALA	ASP	ASP	ARG	ARG	LEU	LEU	LEU	ASP	ASP	THR	THR	THR	THR	PHEN	PHEN	PHEN	ARG
LEU	LEU	LEU	PRO	GLU	ARG	GLU	GLU	VAL	VAL	VAL	VAL	VAL	VAL	VAL	TRP	TRP	TRP	GLN
SER	SER	PRO	PRO	PRO	ALA	ALA	ALA	ALA	VAL	VAL	SER	SER	SER	SER	ASN	ASN	ASN	LYS
LYS	LYS	TRP	TRP	MET	MET	ASP	ASP	SER	VAL	VAL	THR	THR	THR	THR	ILE	ILE	ILE	THR
ASP	ASP	ARG	ARG	ALA	ALA	THR	THR	THR	SER	SER	ASN	ASN	ASN	ASN	ASN	ASN	ASN	LYS
LEU	LEU	ILE	ILE	THR	THR	GLN	GLN	GLN	LYS	LYS	LYS	LYS	LYS	LYS	VAL	VAL	VAL	GLU
PRO	PRO	THR	THR	LEU	LEU	PHEN	PHEN	PHEN	ASP	ASP	LEU	LEU	LEU	LEU	PHEN	PHEN	PHEN	GLY
ASP	ASP	ILE	ASN	SER	SER	LYS	LYS	LYS	SER	SER	LYS	LYS	LYS	LYS	ASP	ASP	ASP	ASP
LEU	LEU	GLU	GLU	GLU	GLY	VAL	VAL	VAL	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	LEU	LEU	LEU	ILE	ILE	ILE	ILE	THR</									

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	90803	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 ($e^-/\text{\AA}^2$)	10	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	2.476	Depositor
Minimum map value	-1.989	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.058	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	315.0, 315.0, 315.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.05, 1.05, 1.05	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.34	0/3097	0.46	0/4247
1	B	0.34	0/3097	0.46	0/4247
All	All	0.34	0/6194	0.46	0/8494

There are no bond length outliers.

There are no bond angle outliers.

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	3031	0	2605	95	0
1	B	3031	0	2605	94	0
All	All	6062	0	5210	189	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 17.

All (189) 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:338:THR:HA	1:A:341:ILE:HD12	1.54	0.89
1:B:338:THR:HA	1:B:341:ILE:HD12	1.54	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LEU:HD12	1:A:533:TYR:HE1	1.49	0.78
1:B:304:LEU:HD12	1:B:533:TYR:HE1	1.49	0.77
1:B:351:ARG:O	1:B:355:LEU:N	2.18	0.75
1:B:339:SER:HB2	1:B:519:PRO:HB3	1.72	0.71
1:A:351:ARG:O	1:A:355:LEU:N	2.18	0.71
1:A:339:SER:HB2	1:A:519:PRO:HB3	1.72	0.71
1:A:744:GLY:HA2	1:A:747:ILE:HD12	1.74	0.69
1:A:448:ASP:OD2	1:A:597:SER:OG	2.12	0.68
1:B:744:GLY:HA2	1:B:747:ILE:HD12	1.74	0.68
1:B:448:ASP:OD2	1:B:597:SER:OG	2.12	0.67
1:A:334:THR:HG21	1:A:686:TYR:HB3	1.81	0.62
1:A:356:ILE:O	1:A:360:LEU:N	2.33	0.62
1:B:356:ILE:O	1:B:360:LEU:N	2.33	0.61
1:B:334:THR:HG21	1:B:686:TYR:HB3	1.81	0.60
1:B:404:ASP:O	1:B:408:ASP:N	2.33	0.60
1:B:313:GLY:O	1:B:546:ARG:NH1	2.35	0.60
1:A:313:GLY:O	1:A:546:ARG:NH1	2.35	0.59
1:A:350:GLY:O	1:A:351:ARG:NE	2.36	0.59
1:B:350:GLY:O	1:B:351:ARG:NE	2.36	0.59
1:A:404:ASP:O	1:A:408:ASP:N	2.33	0.59
1:B:348:ARG:O	1:B:358:ARG:NH2	2.33	0.59
1:B:387:PHE:O	1:B:391:VAL:HG23	2.04	0.58
1:B:456:PRO:HA	1:B:465:PHE:HE1	1.69	0.58
1:A:404:ASP:O	1:A:407:ASN:N	2.34	0.58
1:A:387:PHE:O	1:A:391:VAL:HG23	2.04	0.57
1:B:642:ALA:HB1	1:B:652:LEU:H	1.71	0.56
1:A:670:GLU:O	1:A:673:VAL:HG22	2.05	0.56
1:B:325:MET:HA	1:B:328:THR:HG22	1.88	0.56
1:B:670:GLU:O	1:B:673:VAL:HG22	2.05	0.56
1:A:348:ARG:O	1:A:358:ARG:NH2	2.33	0.56
1:B:404:ASP:O	1:B:407:ASN:N	2.34	0.56
1:A:325:MET:HA	1:A:328:THR:HG22	1.88	0.56
1:A:642:ALA:HB1	1:A:652:LEU:H	1.71	0.56
1:B:344:ASN:OD1	1:B:345:GLY:N	2.39	0.56
1:B:372:PHE:HZ	1:B:503:ALA:HB2	1.70	0.56
1:A:344:ASN:OD1	1:A:345:GLY:N	2.39	0.56
1:A:456:PRO:HA	1:A:465:PHE:HE1	1.69	0.56
1:A:387:PHE:O	1:A:390:THR:HG22	2.07	0.55
1:A:372:PHE:HZ	1:A:503:ALA:HB2	1.70	0.55
1:A:725:ILE:O	1:A:729:VAL:HG23	2.06	0.55
1:B:725:ILE:O	1:B:729:VAL:HG23	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:PHE:O	1:B:390:THR:HG22	2.07	0.53
1:B:336:LEU:O	1:B:339:SER:OG	2.22	0.53
1:A:333:ILE:HG21	1:A:719:GLY:HA3	1.90	0.53
1:B:333:ILE:HG21	1:B:719:GLY:HA3	1.90	0.53
1:B:378:VAL:HG13	1:B:678:ILE:HG13	1.91	0.53
1:B:304:LEU:HD12	1:B:533:TYR:CE1	2.39	0.52
1:A:378:VAL:HG13	1:A:678:ILE:HG13	1.91	0.52
1:B:739:TYR:HA	1:B:742:VAL:HG12	1.92	0.52
1:B:471:GLU:H	1:B:471:GLU:CD	2.18	0.51
1:A:712:ASN:OD1	1:A:714:TRP:N	2.43	0.51
1:A:739:TYR:HA	1:A:742:VAL:HG12	1.92	0.51
1:B:367:ALA:HA	1:B:745:LEU:HD23	1.92	0.51
1:A:367:ALA:HA	1:A:745:LEU:HD23	1.92	0.51
1:B:329:VAL:HA	1:B:332:THR:HG22	1.93	0.51
1:A:329:VAL:HA	1:A:332:THR:HG22	1.93	0.51
1:B:712:ASN:OD1	1:B:714:TRP:N	2.43	0.50
1:B:315:ALA:O	1:B:320:SER:OG	2.29	0.50
1:A:315:ALA:O	1:A:320:SER:OG	2.29	0.50
1:A:471:GLU:H	1:A:471:GLU:CD	2.18	0.50
1:A:743:LEU:O	1:A:747:ILE:HG13	2.12	0.50
1:B:305:PHE:HD1	1:B:593:MET:HG2	1.77	0.49
1:A:693:VAL:HG13	1:A:704:TRP:HZ2	1.77	0.49
1:B:743:LEU:O	1:B:747:ILE:HG13	2.12	0.49
1:A:526:ILE:O	1:A:530:THR:HG23	2.13	0.49
1:B:526:ILE:O	1:B:530:THR:HG23	2.13	0.49
1:B:693:VAL:HG13	1:B:704:TRP:HZ2	1.77	0.49
1:A:550:GLY:HA3	1:A:584:TYR:OH	2.13	0.49
1:A:623:PRO:HG3	1:A:658:THR:HG21	1.95	0.49
1:B:371:ILE:HD11	1:B:681:PHE:HD1	1.78	0.48
1:A:305:PHE:HD1	1:A:593:MET:HG2	1.77	0.48
1:B:727:MET:HG2	1:B:735:ALA:HB2	1.95	0.48
1:A:607:ILE:O	1:A:611:THR:HG23	2.14	0.48
1:B:550:GLY:HA3	1:B:584:TYR:OH	2.13	0.48
1:B:689:ILE:O	1:B:692:SER:OG	2.30	0.48
1:A:371:ILE:HD11	1:A:681:PHE:HD1	1.78	0.47
1:A:304:LEU:HD12	1:A:533:TYR:CE1	2.39	0.47
1:B:623:PRO:HG3	1:B:658:THR:HG21	1.95	0.47
1:A:326:MET:O	1:A:330:VAL:HG23	2.15	0.47
1:B:384:VAL:HG11	1:B:666:ILE:HD11	1.97	0.47
1:B:471:GLU:OE1	1:B:471:GLU:N	2.42	0.47
1:B:607:ILE:O	1:B:611:THR:HG23	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:LEU:O	1:A:339:SER:OG	2.23	0.47
1:B:344:ASN:HB2	1:B:704:TRP:O	2.15	0.47
1:A:316:GLY:H	1:A:479:PRO:HA	1.80	0.47
1:A:344:ASN:HB2	1:A:704:TRP:O	2.15	0.47
1:B:316:GLY:H	1:B:479:PRO:HA	1.80	0.46
1:A:727:MET:HG2	1:A:735:ALA:HB2	1.95	0.46
1:A:438:LEU:HD13	1:A:613:SER:HB3	1.97	0.46
1:A:328:THR:HA	1:A:331:THR:OG1	2.16	0.46
1:A:384:VAL:HG11	1:A:666:ILE:HD11	1.97	0.46
1:A:471:GLU:OE1	1:A:471:GLU:N	2.42	0.46
1:A:636:PRO:HD2	1:A:748:TYR:HE2	1.80	0.46
1:B:636:PRO:HD2	1:B:748:TYR:HE2	1.80	0.46
1:A:302:VAL:O	1:A:306:ILE:HG12	2.16	0.46
1:B:438:LEU:HD13	1:B:613:SER:HB3	1.97	0.46
1:A:701:SER:N	1:A:702:PRO:HD2	2.31	0.46
1:B:326:MET:O	1:B:330:VAL:HG23	2.15	0.46
1:B:496:PRO:O	1:B:499:THR:OG1	2.34	0.46
1:B:328:THR:HA	1:B:331:THR:OG1	2.16	0.46
1:B:701:SER:N	1:B:702:PRO:HD2	2.31	0.46
1:A:372:PHE:CZ	1:A:503:ALA:HB2	2.51	0.45
1:A:689:ILE:O	1:A:692:SER:OG	2.30	0.45
1:B:303:MET:HA	1:B:307:ARG:HB2	1.99	0.45
1:A:303:MET:HA	1:A:307:ARG:HB2	1.99	0.45
1:B:302:VAL:O	1:B:306:ILE:HG12	2.16	0.45
1:B:712:ASN:HD21	1:B:715:ILE:HG23	1.83	0.44
1:B:343:THR:OG1	1:B:707:ALA:HB3	2.17	0.44
1:A:640:MET:O	1:A:653:ARG:NE	2.51	0.44
1:B:324:ILE:HD13	1:B:533:TYR:HB3	2.01	0.43
1:A:343:THR:OG1	1:A:707:ALA:HB3	2.17	0.43
1:B:307:ARG:O	1:B:310:TRP:HB3	2.19	0.43
1:B:307:ARG:HD3	1:B:307:ARG:HA	1.76	0.43
1:B:412:ILE:O	1:B:416:THR:OG1	2.30	0.43
1:B:640:MET:O	1:B:653:ARG:NE	2.51	0.43
1:B:718:LEU:HD12	1:B:722:LEU:HD23	2.00	0.43
1:A:308:LEU:HD12	1:A:308:LEU:HA	1.76	0.43
1:A:501:ILE:HG13	1:A:502:LEU:H	1.83	0.43
1:A:686:TYR:HA	1:A:689:ILE:HG22	2.01	0.43
1:B:654:GLY:O	1:B:658:THR:HG22	2.19	0.43
1:A:748:TYR:HA	1:A:751:TYR:CE1	2.54	0.43
1:A:324:ILE:HD13	1:A:533:TYR:HB3	2.01	0.43
1:A:371:ILE:HD11	1:A:681:PHE:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:PHE:O	1:A:453:THR:HG23	2.19	0.43
1:B:622:ALA:HB3	1:B:623:PRO:HD3	2.01	0.43
1:A:622:ALA:HB3	1:A:623:PRO:HD3	2.01	0.43
1:B:373:ALA:HB1	1:B:626:PHE:CE2	2.54	0.43
1:B:449:PHE:O	1:B:453:THR:HG23	2.19	0.43
1:A:419:ILE:O	1:A:423:ILE:HG12	2.19	0.43
1:A:654:GLY:O	1:A:658:THR:HG22	2.19	0.43
1:B:382:MET:O	1:B:385:VAL:HG12	2.19	0.43
1:B:748:TYR:HA	1:B:751:TYR:CE1	2.54	0.43
1:A:712:ASN:HD21	1:A:715:ILE:HG23	1.83	0.42
1:B:297:LEU:HD23	1:B:297:LEU:HA	1.79	0.42
1:B:372:PHE:CZ	1:B:503:ALA:HB2	2.51	0.42
1:A:486:THR:N	1:A:489:SER:OG	2.52	0.42
1:A:690:ASN:OD1	1:A:691:PHE:N	2.53	0.42
1:B:501:ILE:HG13	1:B:502:LEU:H	1.83	0.42
1:B:670:GLU:O	1:B:674:ILE:HD12	2.19	0.42
1:B:686:TYR:HA	1:B:689:ILE:HG22	2.01	0.42
1:A:495:PHE:N	1:A:496:PRO:HD2	2.34	0.42
1:A:718:LEU:HD12	1:A:722:LEU:HD23	2.00	0.42
1:B:419:ILE:O	1:B:423:ILE:HG12	2.19	0.42
1:B:311:ILE:HG12	1:B:490:VAL:HG13	2.01	0.42
1:B:476:ASN:ND2	1:B:541:GLY:O	2.39	0.42
1:B:545:VAL:HG22	1:B:546:ARG:H	1.85	0.42
1:B:572:PHE:O	1:B:575:SER:OG	2.35	0.42
1:B:683:LEU:HA	1:B:683:LEU:HD23	1.85	0.42
1:A:542:SER:O	1:A:542:SER:OG	2.36	0.42
1:B:495:PHE:N	1:B:496:PRO:HD2	2.34	0.42
1:A:307:ARG:O	1:A:310:TRP:HB3	2.19	0.42
1:A:373:ALA:HB1	1:A:626:PHE:CE2	2.54	0.42
1:B:381:ALA:HA	1:B:384:VAL:HG12	2.02	0.42
1:A:412:ILE:O	1:A:416:THR:OG1	2.30	0.42
1:A:545:VAL:HG22	1:A:546:ARG:H	1.85	0.42
1:B:486:THR:N	1:B:489:SER:OG	2.52	0.42
1:A:476:ASN:ND2	1:A:541:GLY:O	2.39	0.42
1:A:496:PRO:O	1:A:499:THR:OG1	2.34	0.42
1:A:585:GLY:C	1:A:587:MET:H	2.28	0.42
1:A:334:THR:O	1:A:338:THR:HG22	2.20	0.42
1:A:339:SER:O	1:A:343:THR:HG22	2.20	0.42
1:A:381:ALA:HA	1:A:384:VAL:HG12	2.02	0.42
1:A:382:MET:O	1:A:385:VAL:HG12	2.19	0.42
1:A:653:ARG:HE	1:A:653:ARG:HB3	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ILE:HG12	1:A:490:VAL:HG13	2.01	0.41
1:A:670:GLU:O	1:A:674:ILE:HD12	2.19	0.41
1:B:371:ILE:HD11	1:B:681:PHE:CD1	2.54	0.41
1:B:690:ASN:OD1	1:B:691:PHE:N	2.53	0.41
1:B:334:THR:O	1:B:338:THR:HG22	2.20	0.41
1:B:339:SER:O	1:B:343:THR:HG22	2.20	0.41
1:A:305:PHE:HA	1:A:593:MET:HE3	2.02	0.41
1:A:693:VAL:HG13	1:A:704:TRP:CZ2	2.54	0.41
1:B:313:GLY:HA3	1:B:586:LEU:O	2.21	0.41
1:B:388:ALA:O	1:B:392:VAL:HG13	2.21	0.41
1:B:593:MET:HE2	1:B:593:MET:HB2	1.90	0.41
1:A:297:LEU:HD23	1:A:297:LEU:HA	1.79	0.41
1:A:687:ALA:HB2	1:A:723:CYS:HB3	2.03	0.41
1:B:693:VAL:HG13	1:B:704:TRP:CZ2	2.54	0.41
1:A:663:LEU:O	1:A:667:LEU:HG	2.21	0.41
1:B:305:PHE:HA	1:B:593:MET:HE3	2.03	0.40
1:B:663:LEU:O	1:B:667:LEU:HG	2.21	0.40
1:A:313:GLY:HA3	1:A:586:LEU:O	2.21	0.40
1:A:388:ALA:O	1:A:392:VAL:HG13	2.21	0.40
1:B:328:THR:O	1:B:332:THR:HG22	2.21	0.40
1:A:328:THR:O	1:A:332:THR:HG22	2.21	0.40
1:B:687:ALA:HB2	1:B:723:CYS:HB3	2.03	0.40
1:A:745:LEU:HD12	1:A:745:LEU:HA	1.82	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	464/1212 (38%)	407 (88%)	57 (12%)	0	100	100
1	B	464/1212 (38%)	407 (88%)	57 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	928/2424 (38%)	814 (88%)	114 (12%)	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	233/984 (24%)	232 (100%)	1 (0%)	84	81
1	B	233/984 (24%)	232 (100%)	1 (0%)	84	81
All	All	466/1968 (24%)	464 (100%)	2 (0%)	81	81

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	740	VAL
1	A	740	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	298	ASN
1	B	398	HIS
1	A	298	ASN
1	A	398	HIS

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

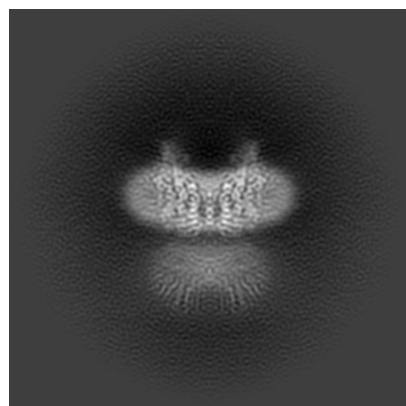
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20537. These allow visual inspection of the internal detail of the map and identification of artifacts.

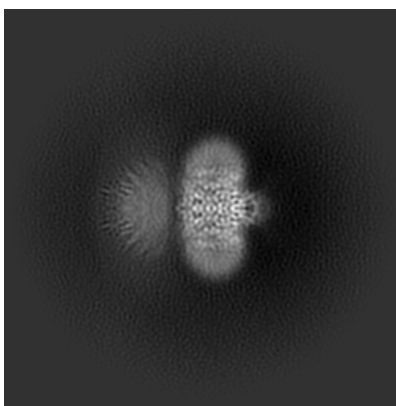
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

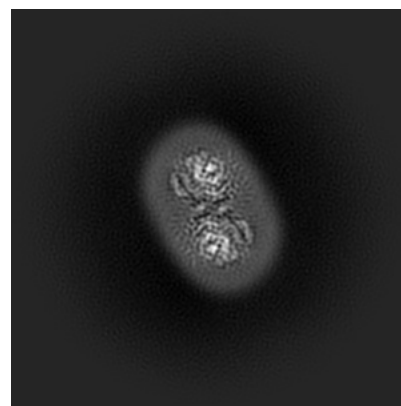
6.1.1 Primary map



X

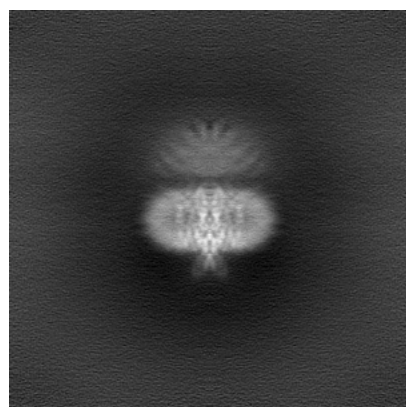


Y

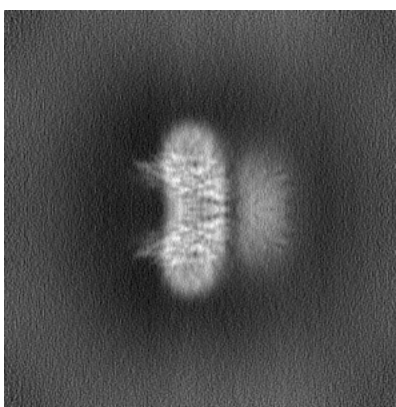


Z

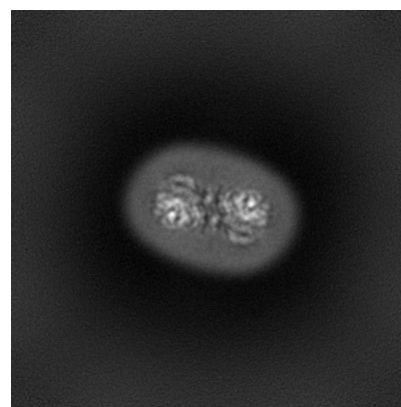
6.1.2 Raw map



X



Y

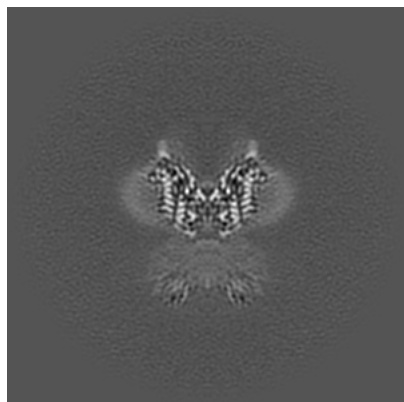


Z

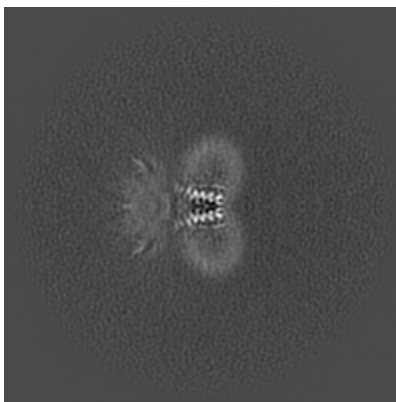
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

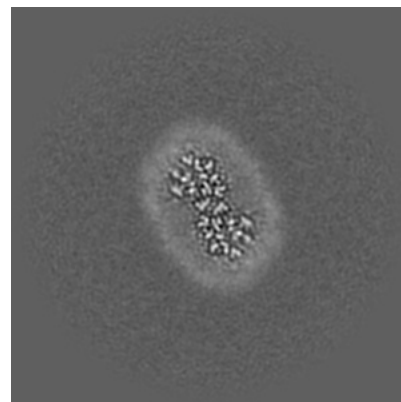
6.2.1 Primary map



X Index: 150

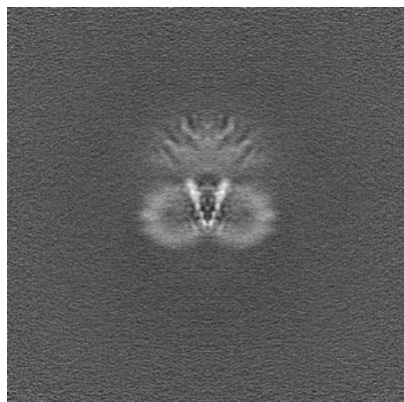


Y Index: 150

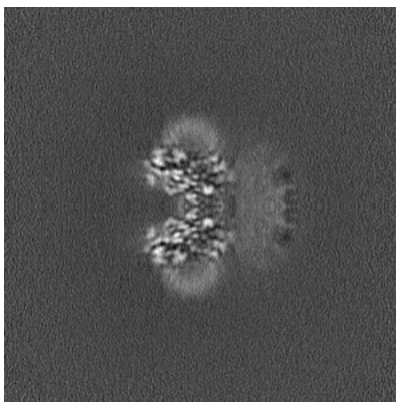


Z Index: 150

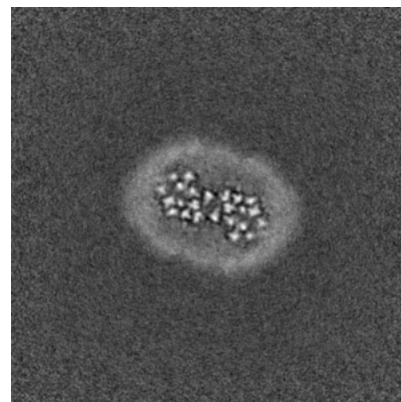
6.2.2 Raw map



X Index: 150



Y Index: 150

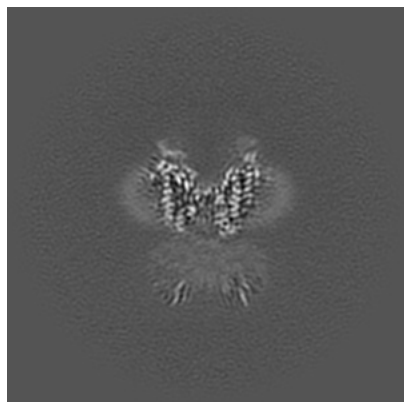


Z Index: 150

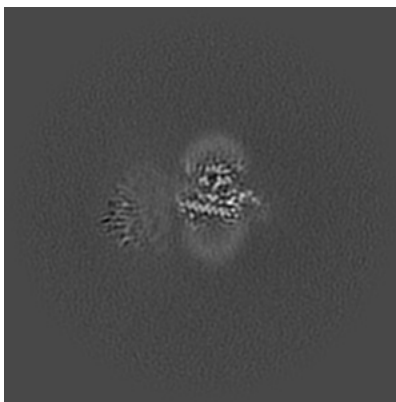
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

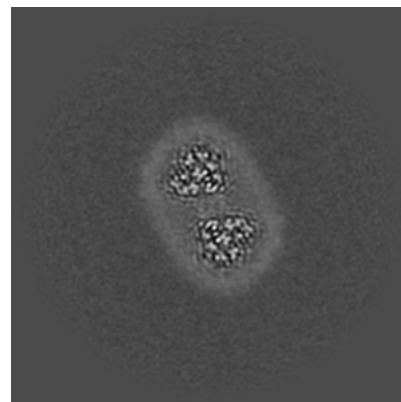
6.3.1 Primary map



X Index: 152

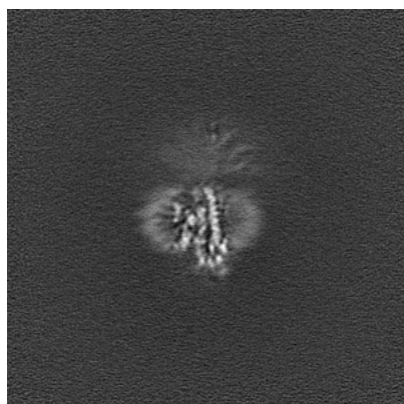


Y Index: 129

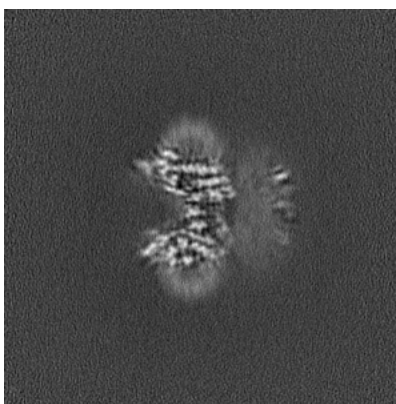


Z Index: 167

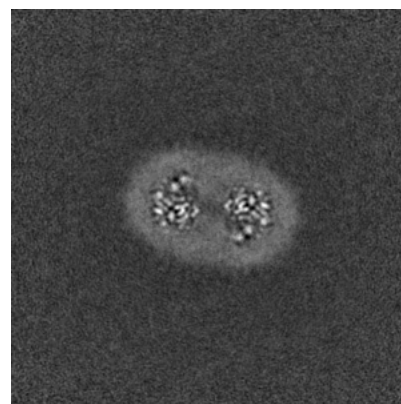
6.3.2 Raw map



X Index: 179



Y Index: 154

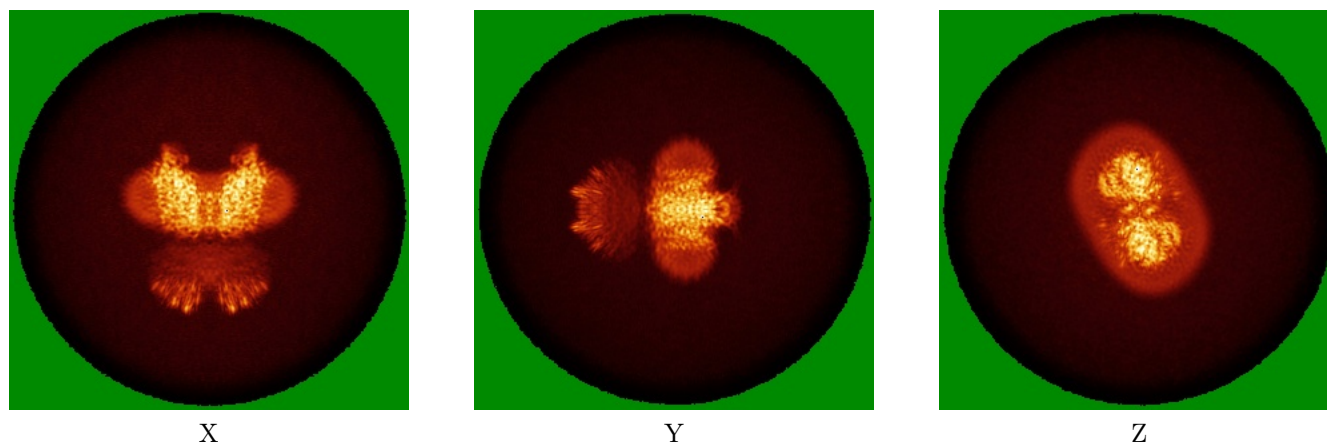


Z Index: 128

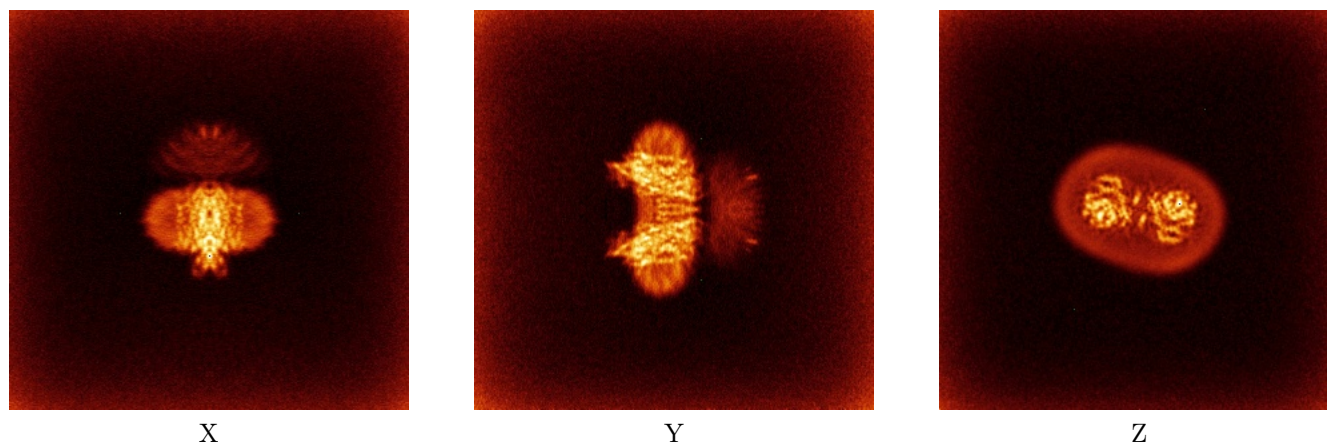
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

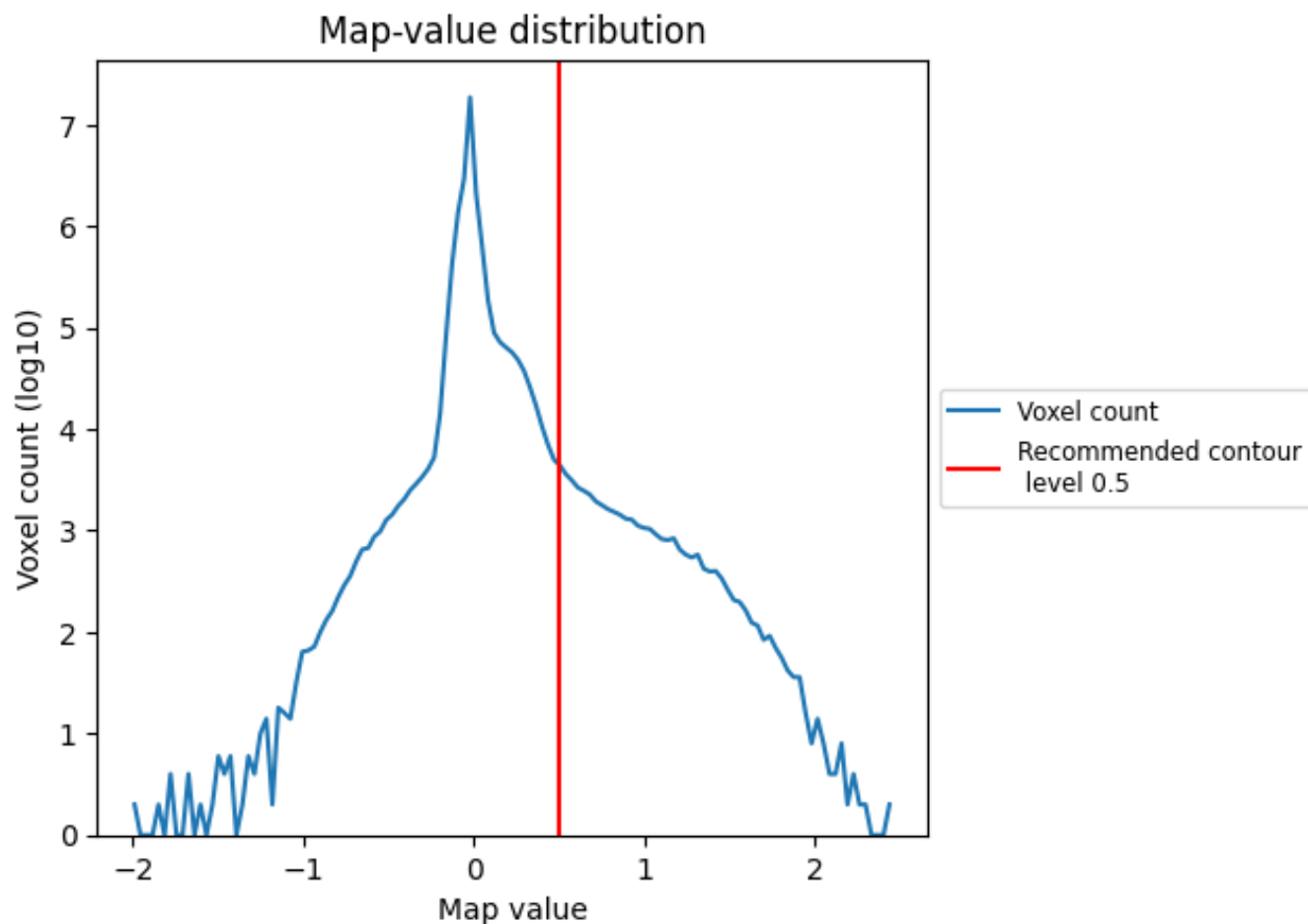
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

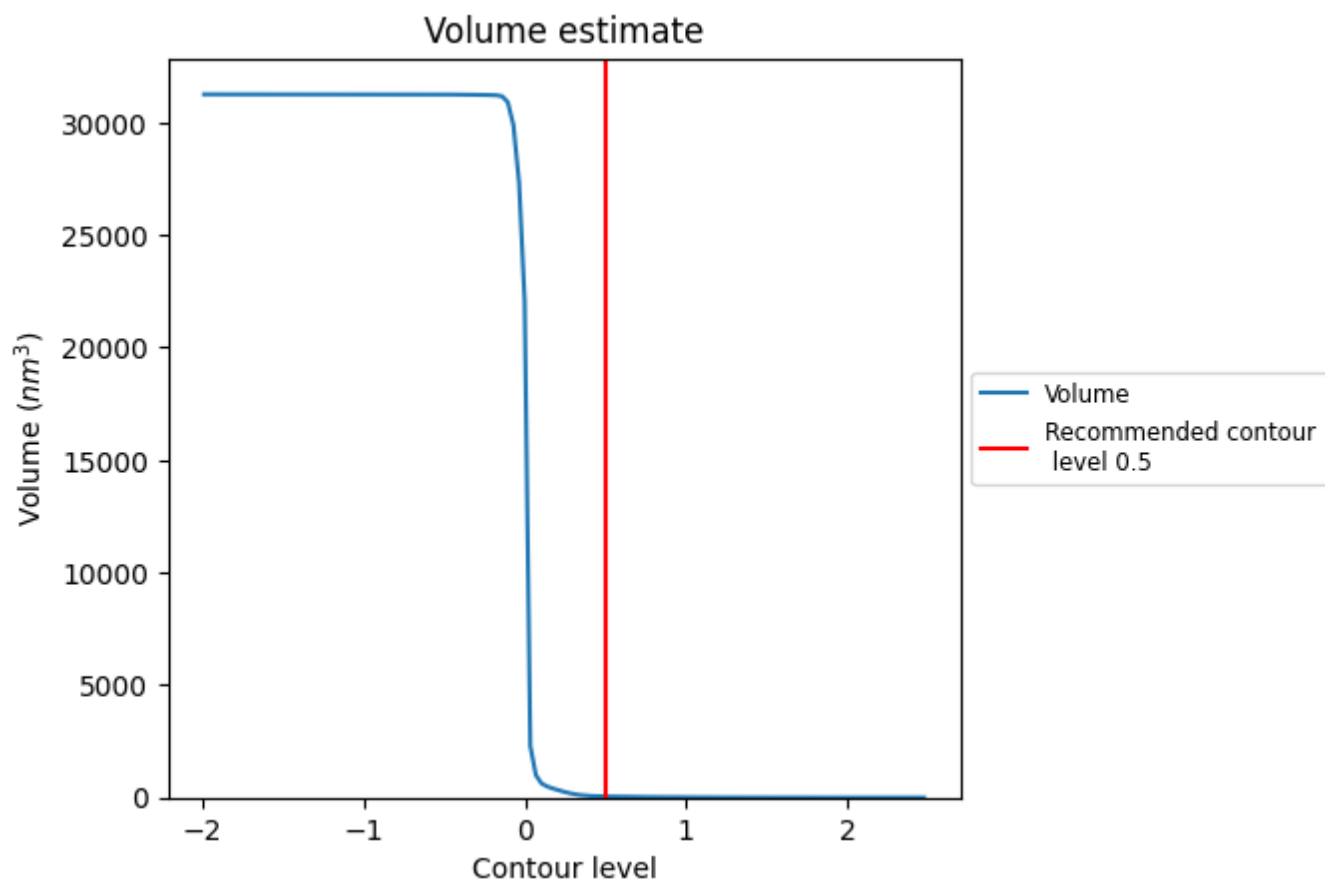
This section contains the results of statistical analysis of the map.

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.

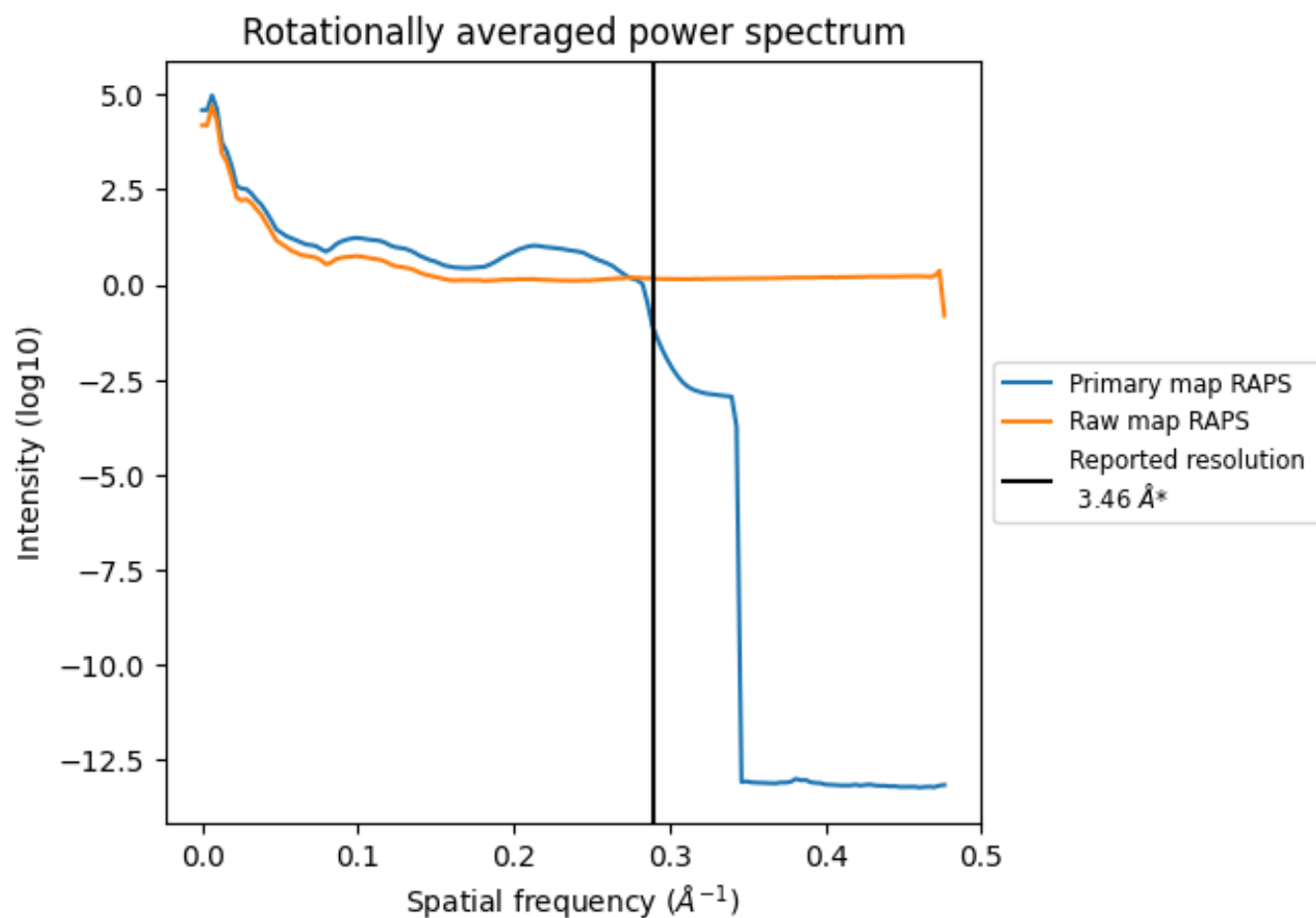
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 49 nm³; this corresponds to an approximate mass of 44 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 ⓘ

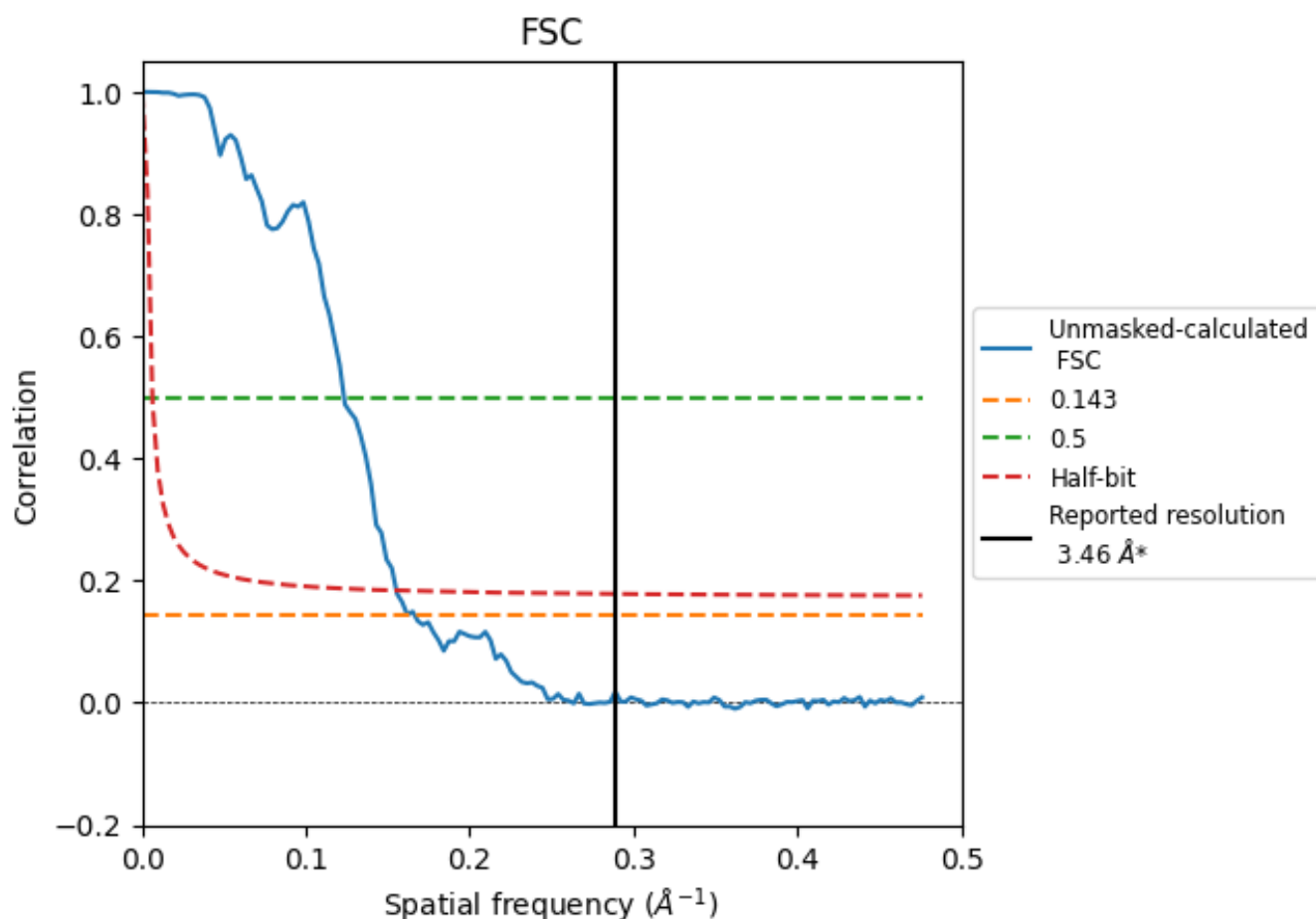


*Reported resolution corresponds to spatial frequency of 0.289 Å⁻¹

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 0.289 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

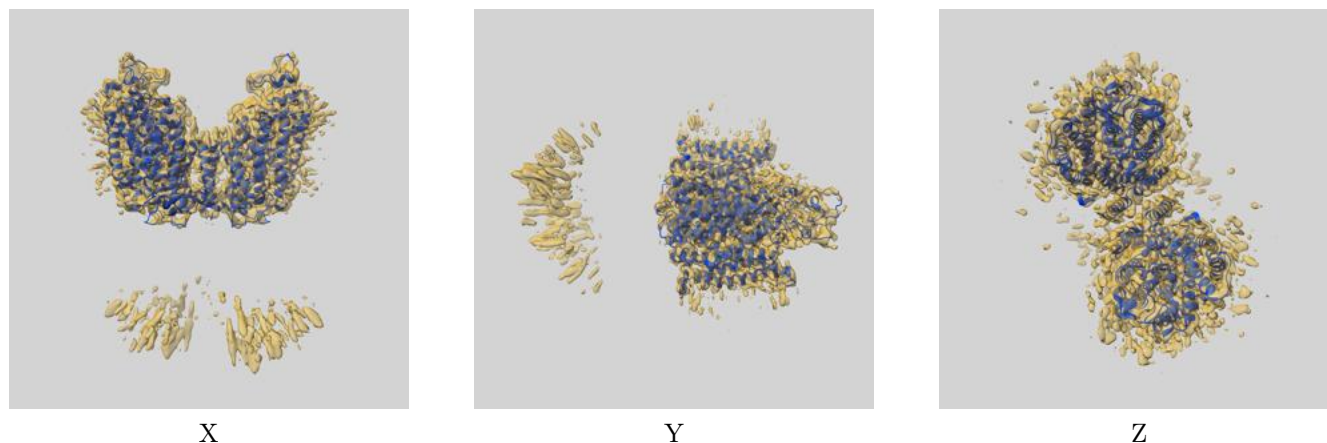
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.46	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.01	8.12	6.44

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.01 differs from the reported value 3.46 by more than 10 %

9 Map-model fit [i](#)

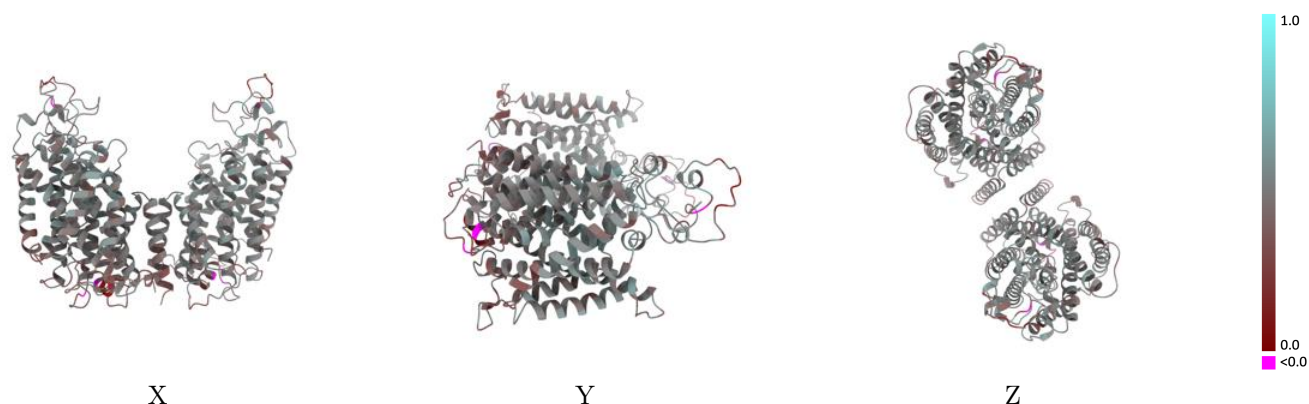
This section contains information regarding the fit between EMDB map EMD-20537 and PDB model 6PZT. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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

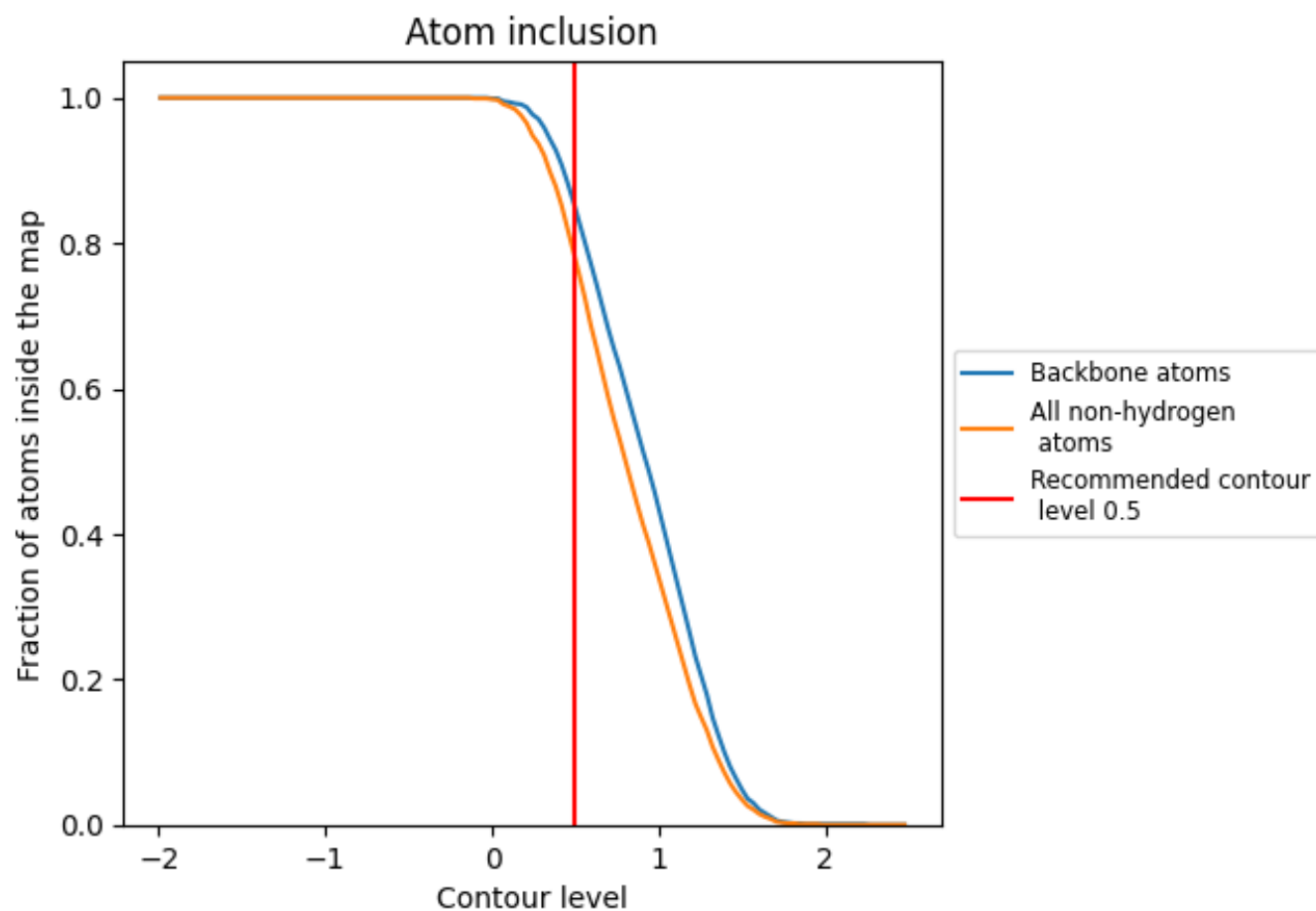


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 78% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7780	0.4540
A	0.7780	0.4550
B	0.7780	0.4540

