



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 09:56 PM UTC

PDB ID : 7UKD / pdb_00007ukd
EMDB ID : EMD-26577
Title : Human Kv4.2-KChIP2 channel complex in an inactivated state, class 2, trans-membrane region
Authors : Zhao, H.; Dai, Y.; Lee, C.H.
Deposited on : 2022-04-01
Resolution : 2.88 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

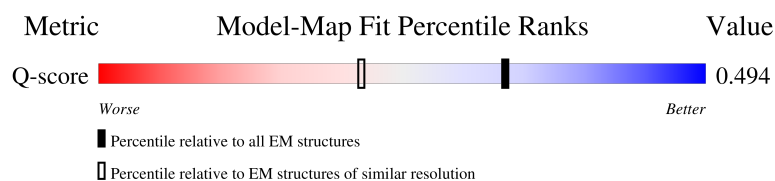
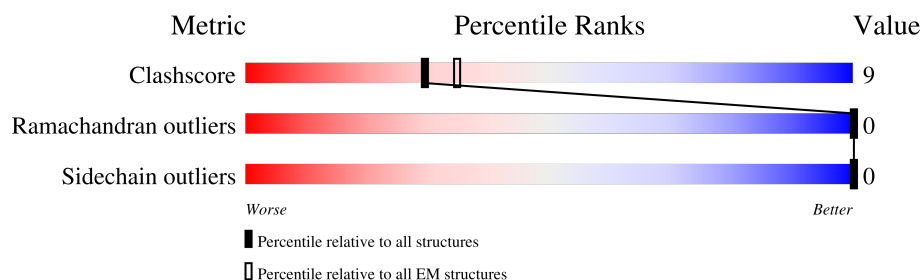
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.88 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	12111 (2.38 - 3.38)

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	524	12% 36% 13% 51%
1	B	524	11% 40% 10% 51%
1	C	524	12% 36% 13% 51%
1	D	524	11% 40% 9% 51%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8000 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.

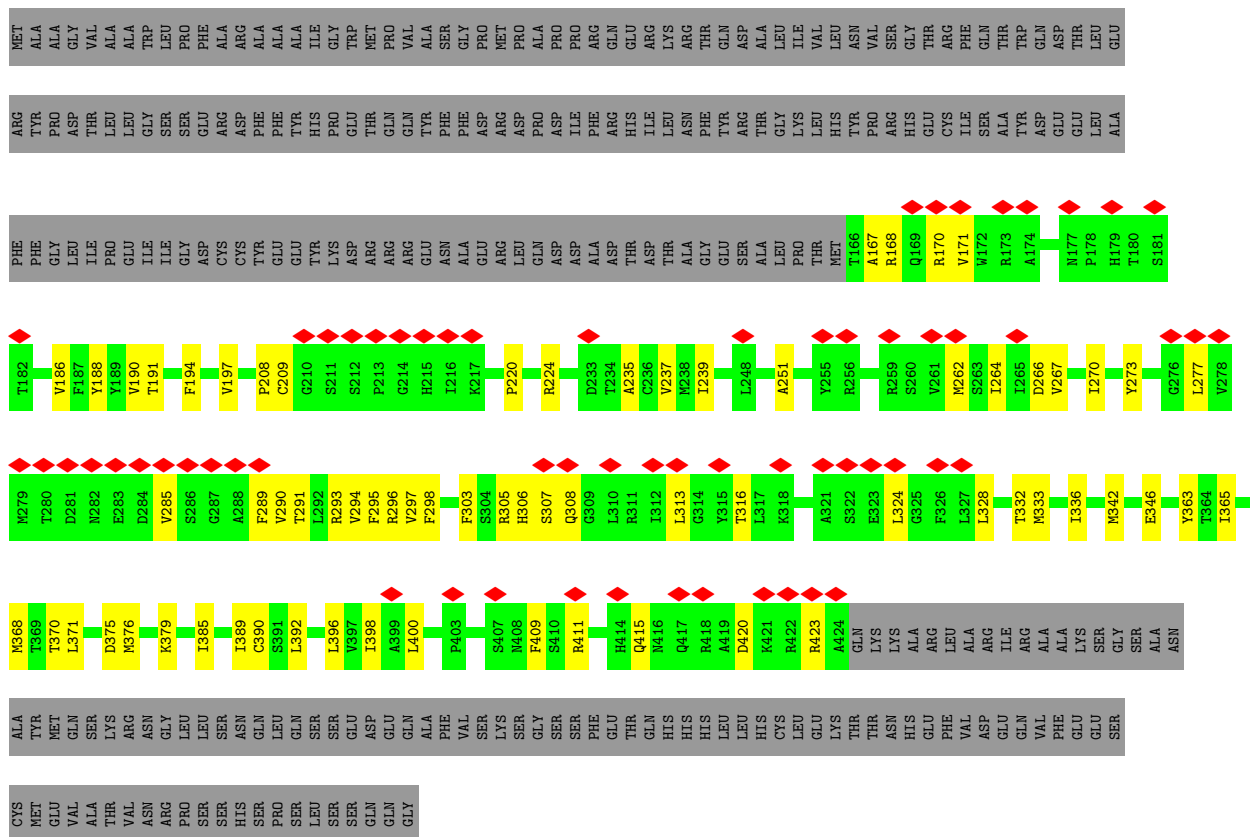
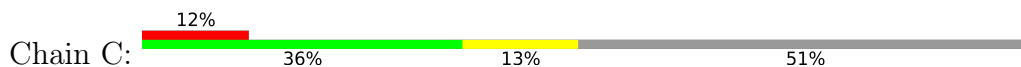
- Molecule 1 is a protein called Potassium voltage-gated channel subfamily D member 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	259	Total 1999	C 1313	N 333	O 339	S 14	0	0
1	B	259	Total 1999	C 1313	N 333	O 339	S 14	0	0
1	C	259	Total 1999	C 1313	N 333	O 339	S 14	0	0
1	D	259	Total 1999	C 1313	N 333	O 339	S 14	0	0

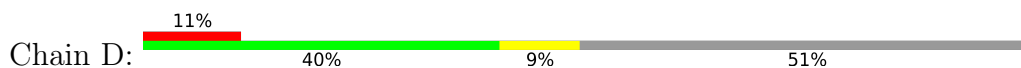
- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
2	A	4	Total 4	K 4	0

- Molecule 1: Potassium voltage-gated channel subfamily D member 2



- Molecule 1: Potassium voltage-gated channel subfamily D member 2



HIS	LEU	HIS	CYS	LEU	GLU	LYS	THR	THR	ASN	HIS	GLU	PHE	VAL	ASP	GLN	GLY	SER	ALA	GLY	SER	ALA	ASN	ARG	PRO	SER	HIS	LEU	SER	SER	GLN	GLN	GLY																					
D420	K421	R422	R423	A424	GLN	LYS	LYS	LYS	ALA	ALA	ALA	LYS	SER	GLY	SER	ALA	ASN	ALA	TYR	MET	VAL	THR	VAL	THR	VAL	VAL	GLY	SER	GLY	GLY	SER	HIS																					
K302	R305	H306	S307	G308	G309	L310	R311	I312	L313	C320	L324	G325	F326	F329	I335	F338	V341	M342	E346	Y363	T364	I365	T370	L371	D375	M376	G383	K384	I385	I389	L392	S393	G394	I398	R411	I412	Y413	H414	Q415	N416	Q417	R418	A419										
Y188	F194	N202	G210	S211	S212	P213	G214	H215	I216	K217	V237	M238	I239	A249	R254	Y255	R266	F267	V268	R269	S260	S263	I264	I265	D266	V267	I270	L271	P272	Y273	L277	V278	M279	T280	D281	N282	E283	D284	V285	S286	G287	A288	F289	R293	V294	F295	V297						
PHE	PHE	GLY	LEU	ILE	PRO	GLU	ILE	ILE	GLY	ASP	CYS	CYS	TYR	GLU	GLU	TYR	LYS	ASP	ARG	ARG	ARG	ASN	ALA	GLU	ARG	ARG	LEU	GLN	ASP	ASP	GLY	ALA	ALA	LEU	PRO	THR	MET	T166	A167	R168	Q169	R170	R173	N177	P178	H179	T180	S181	T182				
ARG	TYR	PRO	ASP	THR	VAL	LEU	LEU	GLY	SER	PRO	PHE	GLU	ASP	ARG	PHE	GLN	TYR	ILE	GLY	THR	THR	GLN	VAL	ALA	TYR	GLY	LEU	ILE	VAL	LEU	HIS	TYR	PRO	VAL	ASN	LEU	THR	GLY	THR	ARG	ILE	PHE	GLN	ALA	THR	TRP	GLN	ASP	GLU	THR	LEU	LEU	ALA

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	210659	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	76.4	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.270	Depositor
Minimum map value	-2.264	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.059	Depositor
Recommended contour level	0.5	Depositor
Map size ($\text{\AA}$)	290.75198, 290.75198, 290.75198	wwPDB
Map dimensions	352, 352, 352	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.826, 0.826, 0.826	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: K

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	0.19	0/2047	0.38	0/2781
1	B	0.19	0/2047	0.36	0/2781
1	C	0.19	0/2047	0.38	0/2781
1	D	0.19	0/2047	0.36	0/2781
All	All	0.19	0/8188	0.37	0/11124

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	1999	0	2031	48	0
1	B	1999	0	2032	37	0
1	C	1999	0	2031	48	0
1	D	1999	0	2032	34	0
2	A	4	0	0	0	0
All	All	8000	0	8126	149	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 9.

All (149) 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:262:MET:HE3	1:A:305:ARG:HB2	1.60	0.83
1:C:262:MET:HE3	1:C:305:ARG:HB2	1.60	0.83
1:C:370:THR:HA	1:D:371:LEU:HD23	1.71	0.71
1:B:320:CYS:SG	1:B:408:ASN:ND2	2.62	0.70
1:A:370:THR:HA	1:B:371:LEU:HD23	1.72	0.70
1:B:173:ARG:HH22	1:B:180:THR:HG21	1.59	0.67
1:D:173:ARG:HH22	1:D:180:THR:HG21	1.59	0.67
1:A:371:LEU:HD23	1:D:370:THR:HA	1.80	0.64
1:B:329:PHE:HZ	1:C:303:PHE:HB3	1.63	0.64
1:C:285:VAL:HG13	1:C:289:PHE:HB2	1.80	0.64
1:A:285:VAL:HG13	1:A:289:PHE:HB2	1.80	0.62
1:D:266:ASP:OD2	1:D:302:LYS:NZ	2.33	0.61
1:B:266:ASP:OD2	1:B:302:LYS:NZ	2.33	0.61
1:C:316:THR:HG21	1:C:409:PHE:HA	1.83	0.61
1:B:370:THR:HA	1:C:371:LEU:HD23	1.83	0.60
1:C:209:CYS:HB3	1:C:224:ARG:HH11	1.68	0.59
1:A:316:THR:HG21	1:A:409:PHE:HA	1.83	0.59
1:A:209:CYS:HB3	1:A:224:ARG:HH11	1.68	0.58
1:A:168:ARG:NH2	1:A:251:ALA:O	2.37	0.57
1:C:168:ARG:NH2	1:C:251:ALA:O	2.37	0.57
1:D:260:SER:O	1:D:263:SER:OG	2.19	0.56
1:A:365:ILE:HD11	1:D:392:LEU:HD21	1.87	0.56
1:B:392:LEU:HD21	1:C:365:ILE:HD11	1.87	0.55
1:B:375:ASP:OD1	1:B:376:MET:N	2.40	0.55
1:B:293:ARG:HA	1:B:296:ARG:HD3	1.89	0.55
1:D:188:TYR:OH	1:D:305:ARG:O	2.24	0.55
1:A:303:PHE:HB3	1:D:329:PHE:HZ	1.71	0.55
1:D:293:ARG:HA	1:D:296:ARG:HD3	1.89	0.55
1:D:375:ASP:OD1	1:D:376:MET:N	2.40	0.55
1:B:188:TYR:OH	1:B:305:ARG:O	2.24	0.54
1:B:320:CYS:HB3	1:B:324:LEU:HG	1.90	0.54
1:A:346:GLU:OE2	1:A:379:LYS:N	2.39	0.53
1:C:167:ALA:HA	1:C:170:ARG:HE	1.74	0.53
1:D:320:CYS:HB3	1:D:324:LEU:HG	1.89	0.52
1:A:375:ASP:OD1	1:A:376:MET:N	2.43	0.52
1:A:208:PRO:HA	1:A:220:PRO:HA	1.92	0.52
1:C:346:GLU:OE2	1:C:379:LYS:N	2.39	0.52
1:B:260:SER:O	1:B:263:SER:OG	2.19	0.51
1:A:237:VAL:HG13	1:A:270:ILE:HG23	1.93	0.51
1:A:167:ALA:HA	1:A:170:ARG:HE	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:VAL:HG13	1:C:270:ILE:HG23	1.93	0.51
1:D:294:VAL:O	1:D:297:VAL:HG12	2.11	0.51
1:B:294:VAL:O	1:B:297:VAL:HG12	2.11	0.50
1:C:208:PRO:HA	1:C:220:PRO:HA	1.93	0.50
1:C:396:LEU:O	1:C:400:LEU:HG	2.11	0.50
1:A:420:ASP:HA	1:A:423:ARG:NE	2.27	0.50
1:A:385:ILE:O	1:A:389:ILE:HG12	2.12	0.50
1:C:375:ASP:OD1	1:C:376:MET:N	2.43	0.50
1:A:396:LEU:O	1:A:400:LEU:HG	2.11	0.49
1:D:237:VAL:HG11	1:D:273:TYR:CD2	2.47	0.49
1:C:420:ASP:HA	1:C:423:ARG:NE	2.27	0.49
1:C:294:VAL:O	1:C:297:VAL:HG12	2.13	0.49
1:A:398:ILE:HD11	1:D:398:ILE:HG13	1.94	0.49
1:B:237:VAL:HG11	1:B:273:TYR:CD2	2.47	0.49
1:C:385:ILE:O	1:C:389:ILE:HG12	2.12	0.49
1:D:394:GLY:O	1:D:398:ILE:HG12	2.13	0.48
1:B:394:GLY:O	1:B:398:ILE:HG12	2.13	0.48
1:C:293:ARG:HD2	1:C:296:ARG:HE	1.79	0.48
1:D:335:ILE:HD11	1:D:365:ILE:HD13	1.95	0.48
1:A:293:ARG:HD2	1:A:296:ARG:HE	1.79	0.48
1:C:303:PHE:O	1:C:307:SER:N	2.31	0.48
1:B:335:ILE:HD11	1:B:365:ILE:HD13	1.95	0.48
1:B:398:ILE:HG13	1:C:398:ILE:HD11	1.96	0.48
1:A:294:VAL:O	1:A:297:VAL:HG12	2.13	0.47
1:C:333:MET:SD	1:D:310:LEU:HD21	2.54	0.47
1:A:235:ALA:O	1:A:239:ILE:HG12	2.14	0.47
1:C:235:ALA:O	1:C:239:ILE:HG12	2.14	0.47
1:A:313:LEU:HD12	1:D:326:PHE:HE2	1.79	0.47
1:B:267:VAL:HA	1:B:270:ILE:HG22	1.96	0.47
1:B:385:ILE:O	1:B:389:ILE:HG12	2.13	0.47
1:A:266:ASP:O	1:A:270:ILE:HD12	2.14	0.47
1:D:385:ILE:O	1:D:389:ILE:HG12	2.14	0.47
1:A:295:PHE:HD1	1:A:298:PHE:HE1	1.63	0.47
1:A:411:ARG:O	1:A:415:GLN:HG2	2.15	0.47
1:C:266:ASP:O	1:C:270:ILE:HD12	2.14	0.47
1:B:194:PHE:CE2	1:B:239:ILE:HG13	2.50	0.47
1:C:411:ARG:O	1:C:415:GLN:HG2	2.15	0.47
1:D:194:PHE:CE2	1:D:239:ILE:HG13	2.50	0.46
1:A:273:TYR:CE2	1:A:277:LEU:HD11	2.51	0.46
1:C:273:TYR:CE2	1:C:277:LEU:HD11	2.51	0.46
1:D:267:VAL:HA	1:D:270:ILE:HG22	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:PHE:O	1:A:307:SER:N	2.31	0.45
1:A:264:ILE:HA	1:A:267:VAL:HG12	1.98	0.45
1:C:290:VAL:O	1:C:294:VAL:HG23	2.16	0.45
1:C:295:PHE:HD1	1:C:298:PHE:HE1	1.63	0.45
1:A:290:VAL:O	1:A:294:VAL:HG23	2.16	0.45
1:D:188:TYR:OH	1:D:306:HIS:HA	2.17	0.45
1:B:188:TYR:OH	1:B:306:HIS:HA	2.17	0.45
1:C:264:ILE:HA	1:C:267:VAL:HG12	1.98	0.45
1:C:295:PHE:HD1	1:C:298:PHE:CE1	2.35	0.45
1:C:324:LEU:O	1:C:328:LEU:HD23	2.17	0.45
1:C:194:PHE:HA	1:C:197:VAL:HG12	1.99	0.44
1:A:194:PHE:HA	1:A:197:VAL:HG12	1.99	0.44
1:B:286:SER:HA	1:B:289:PHE:HE1	1.83	0.44
1:B:326:PHE:HE2	1:C:313:LEU:HD12	1.83	0.44
1:A:295:PHE:HD1	1:A:298:PHE:CE1	2.35	0.44
1:C:188:TYR:OH	1:C:306:HIS:O	2.28	0.43
1:C:191:THR:HG23	1:C:306:HIS:NE2	2.33	0.43
1:A:324:LEU:O	1:A:328:LEU:HD23	2.17	0.43
1:D:260:SER:O	1:D:264:ILE:HD12	2.19	0.43
1:D:286:SER:HA	1:D:289:PHE:HE1	1.83	0.43
1:C:420:ASP:O	1:C:423:ARG:HG2	2.18	0.43
1:A:333:MET:SD	1:B:310:LEU:HD21	2.59	0.43
1:A:420:ASP:O	1:A:423:ARG:HG2	2.18	0.43
1:B:338:PHE:HA	1:B:341:VAL:HG22	2.00	0.42
1:D:338:PHE:HA	1:D:341:VAL:HG22	2.00	0.42
1:A:188:TYR:OH	1:A:306:HIS:O	2.28	0.42
1:A:191:THR:HG23	1:A:306:HIS:NE2	2.33	0.42
1:B:307:SER:HB3	1:B:310:LEU:HD13	2.00	0.42
1:B:260:SER:O	1:B:264:ILE:HD12	2.19	0.42
1:C:342:MET:HE2	1:C:363:TYR:CE2	2.54	0.42
1:D:194:PHE:CD2	1:D:239:ILE:HG13	2.54	0.42
1:A:308:GLN:O	1:A:308:GLN:HG3	2.19	0.42
1:A:332:THR:O	1:A:336:ILE:HG13	2.19	0.42
1:D:307:SER:HB3	1:D:310:LEU:HD13	2.00	0.42
1:A:267:VAL:O	1:A:271:LEU:HD23	2.20	0.42
1:C:392:LEU:HD21	1:D:365:ILE:HD11	2.00	0.42
1:D:346:GLU:OE2	1:D:383:GLY:HA3	2.20	0.42
1:B:194:PHE:CD2	1:B:239:ILE:HG13	2.54	0.42
1:C:308:GLN:HG3	1:C:308:GLN:O	2.19	0.42
1:A:209:CYS:N	1:A:219:LEU:O	2.43	0.42
1:A:342:MET:HE2	1:A:363:TYR:CE2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:342:MET:HE2	1:D:363:TYR:CE2	2.55	0.42
1:B:289:PHE:O	1:B:293:ARG:HG3	2.21	0.41
1:C:295:PHE:CD1	1:C:298:PHE:HE1	2.38	0.41
1:C:332:THR:O	1:C:336:ILE:HG13	2.19	0.41
1:A:289:PHE:O	1:A:293:ARG:HG2	2.19	0.41
1:A:313:LEU:HD12	1:D:326:PHE:CE2	2.53	0.41
1:B:270:ILE:HG23	1:B:271:LEU:HD12	2.02	0.41
1:A:186:VAL:O	1:A:190:VAL:HG22	2.21	0.41
1:A:368:MET:HG3	1:A:390:CYS:SG	2.60	0.41
1:B:326:PHE:CE2	1:C:313:LEU:HD12	2.56	0.41
1:C:291:THR:HG22	1:C:295:PHE:CE2	2.55	0.41
1:D:263:SER:O	1:D:267:VAL:HG23	2.20	0.41
1:C:289:PHE:O	1:C:293:ARG:HG2	2.20	0.41
1:B:342:MET:HE2	1:B:363:TYR:CE2	2.55	0.41
1:B:346:GLU:OE2	1:B:383:GLY:HA3	2.20	0.41
1:C:368:MET:HG3	1:C:390:CYS:SG	2.60	0.41
1:A:291:THR:HG22	1:A:295:PHE:CE2	2.55	0.41
1:C:186:VAL:O	1:C:190:VAL:HG22	2.21	0.41
1:D:202:ASN:OD1	1:D:296:ARG:NH2	2.54	0.41
1:B:342:MET:HE1	1:B:378:PRO:HB3	2.03	0.41
1:D:270:ILE:HG23	1:D:271:LEU:HD12	2.02	0.41
1:A:295:PHE:CD1	1:A:298:PHE:HE1	2.38	0.40
1:B:335:ILE:HA	1:B:364:THR:HG21	2.03	0.40
1:B:263:SER:O	1:B:267:VAL:HG23	2.20	0.40
1:A:167:ALA:O	1:A:171:VAL:HG22	2.22	0.40
1:C:167:ALA:O	1:C:171:VAL:HG22	2.22	0.40
1:B:327:LEU:HD21	1:C:313:LEU:HD11	2.03	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	257/524 (49%)	251 (98%)	6 (2%)	0	100	100
1	B	257/524 (49%)	253 (98%)	4 (2%)	0	100	100
1	C	257/524 (49%)	251 (98%)	6 (2%)	0	100	100
1	D	257/524 (49%)	253 (98%)	4 (2%)	0	100	100
All	All	1028/2096 (49%)	1008 (98%)	20 (2%)	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	B	165/447 (37%)	165 (100%)	0	100	100
1	C	211/447 (47%)	211 (100%)	0	100	100
1	D	211/447 (47%)	211 (100%)	0	100	100
All	All	587/1341 (44%)	587 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	B	408	ASN
1	C	179	HIS
1	D	408	ASN

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 4 ligands modelled in this entry, 4 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.

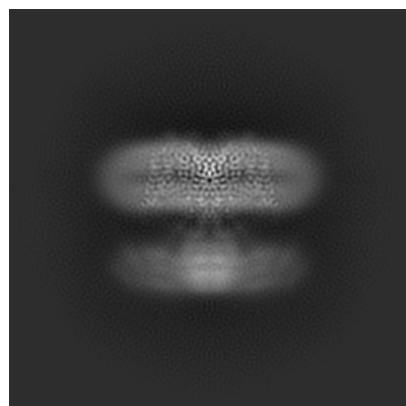
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26577. These allow visual inspection of the internal detail of the map and identification of artifacts.

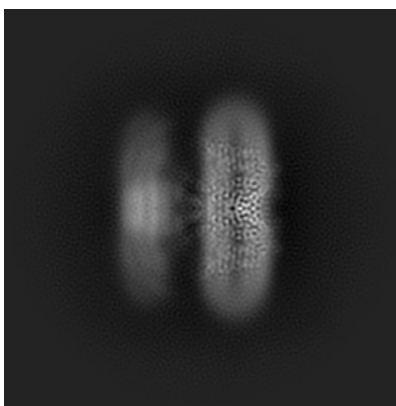
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

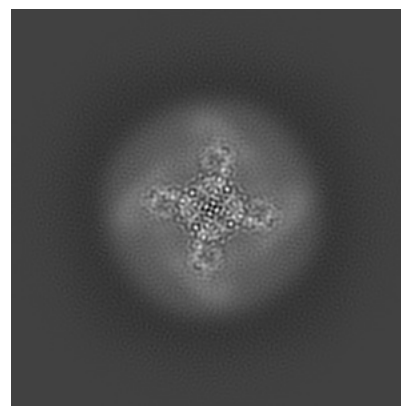
6.1.1 Primary map



X

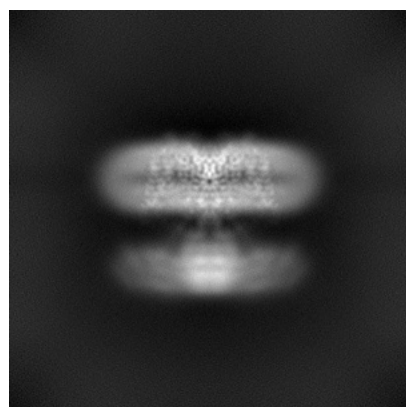


Y

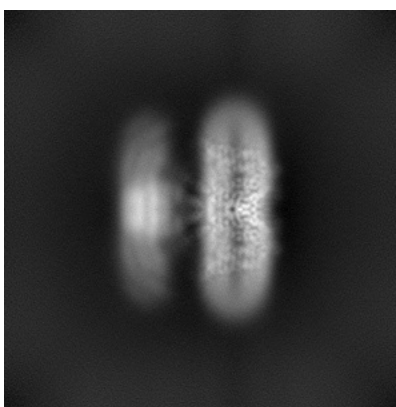


Z

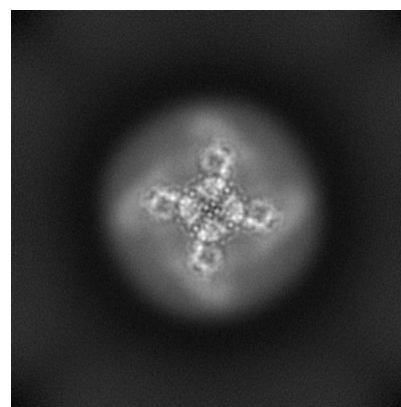
6.1.2 Raw map



X



Y

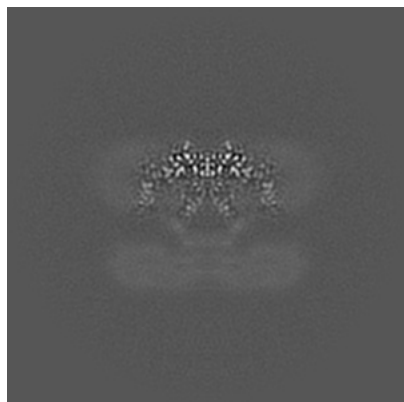


Z

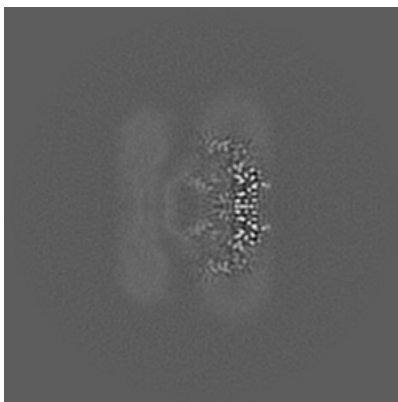
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

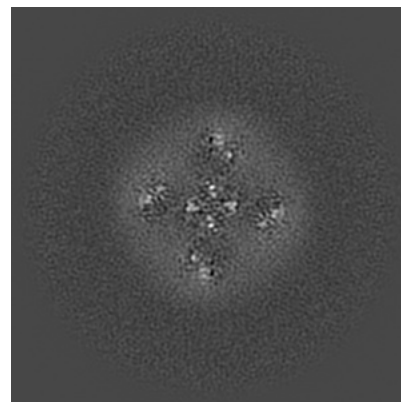
6.2.1 Primary map



X Index: 176

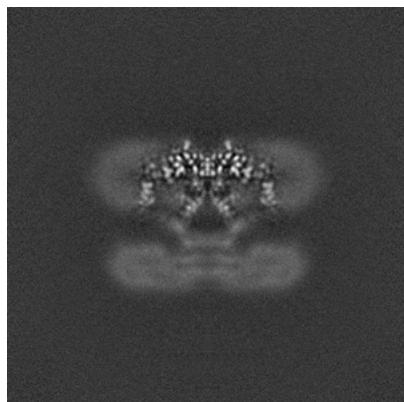


Y Index: 176

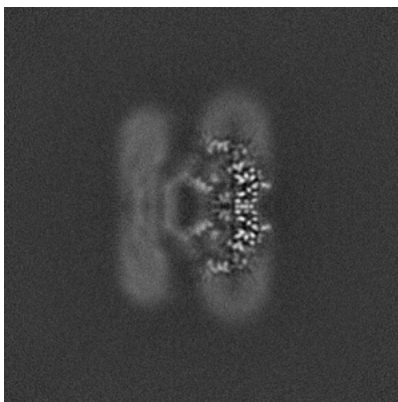


Z Index: 176

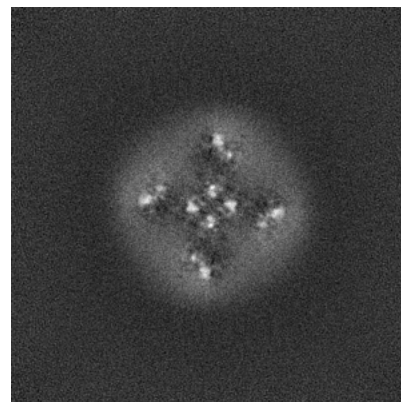
6.2.2 Raw map



X Index: 176



Y Index: 176

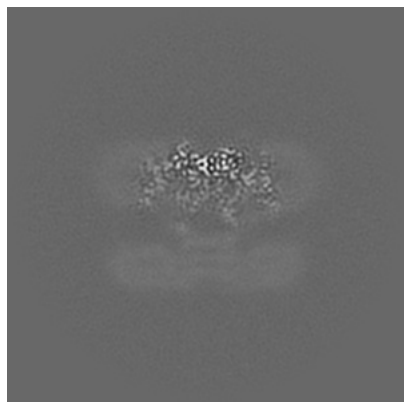


Z Index: 176

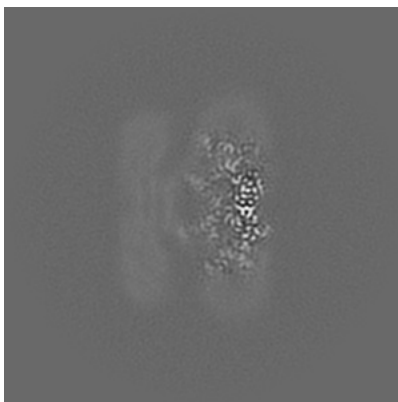
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

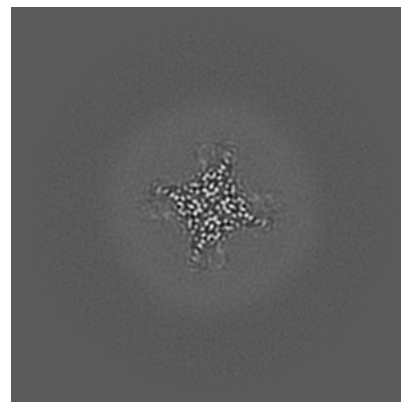
6.3.1 Primary map



X Index: 180

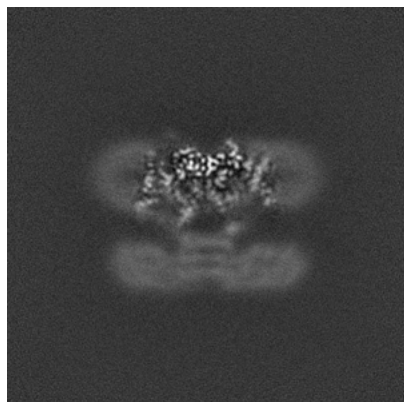


Y Index: 172

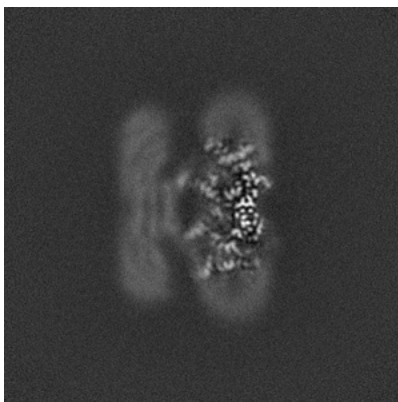


Z Index: 220

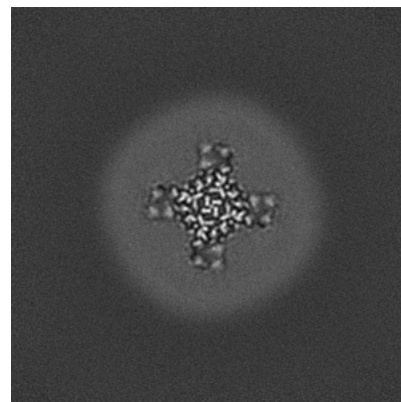
6.3.2 Raw map



X Index: 172



Y Index: 180

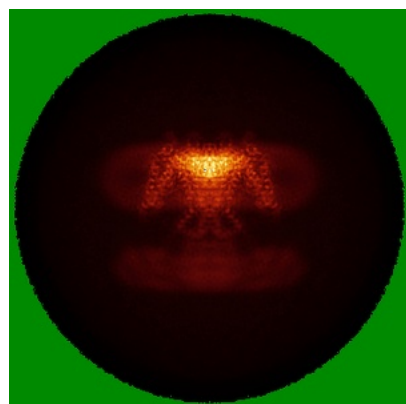


Z Index: 216

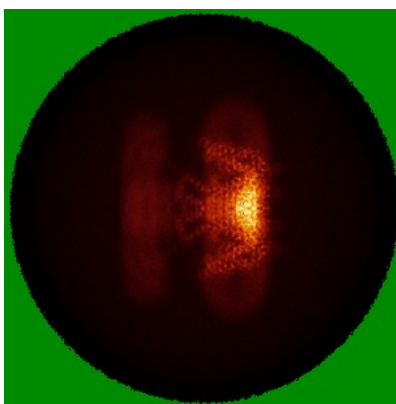
The images above show the largest variance slices of the map in three orthogonal directions.

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

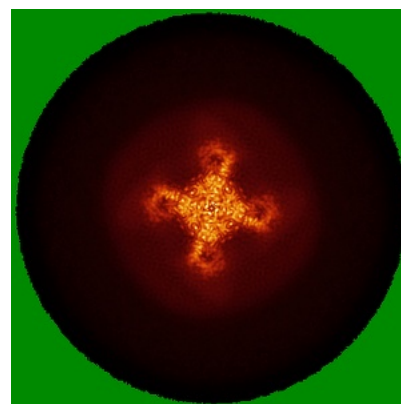
6.4.1 Primary map



X

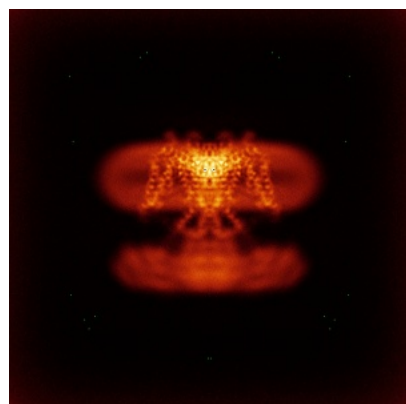


Y

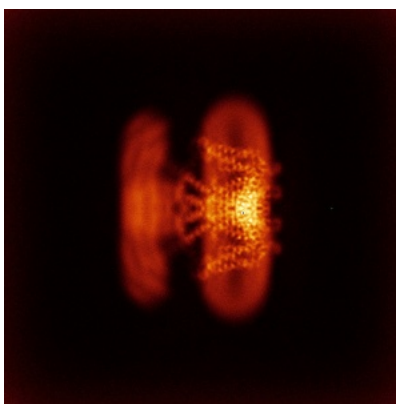


Z

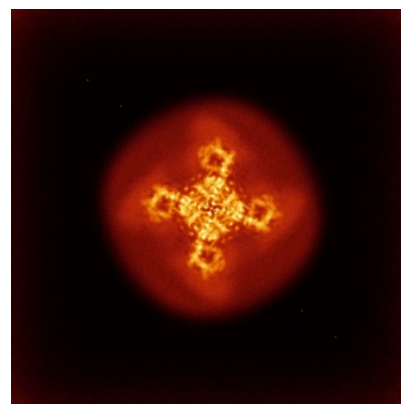
6.4.2 Raw map



X



Y

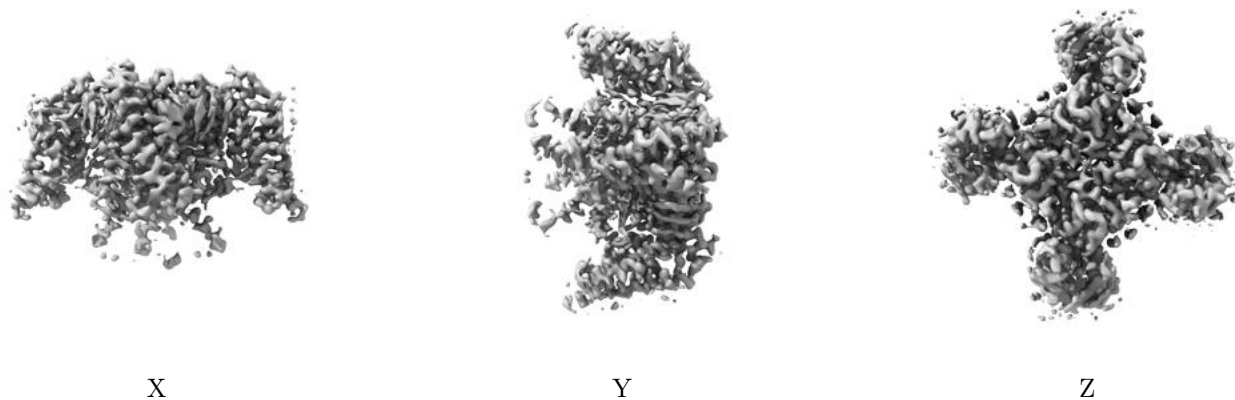


Z

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.

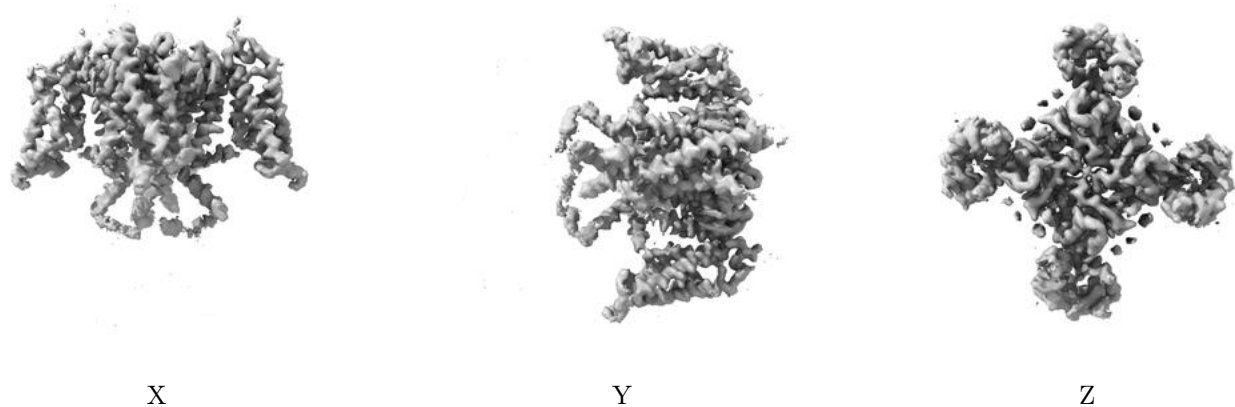
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.5. 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.

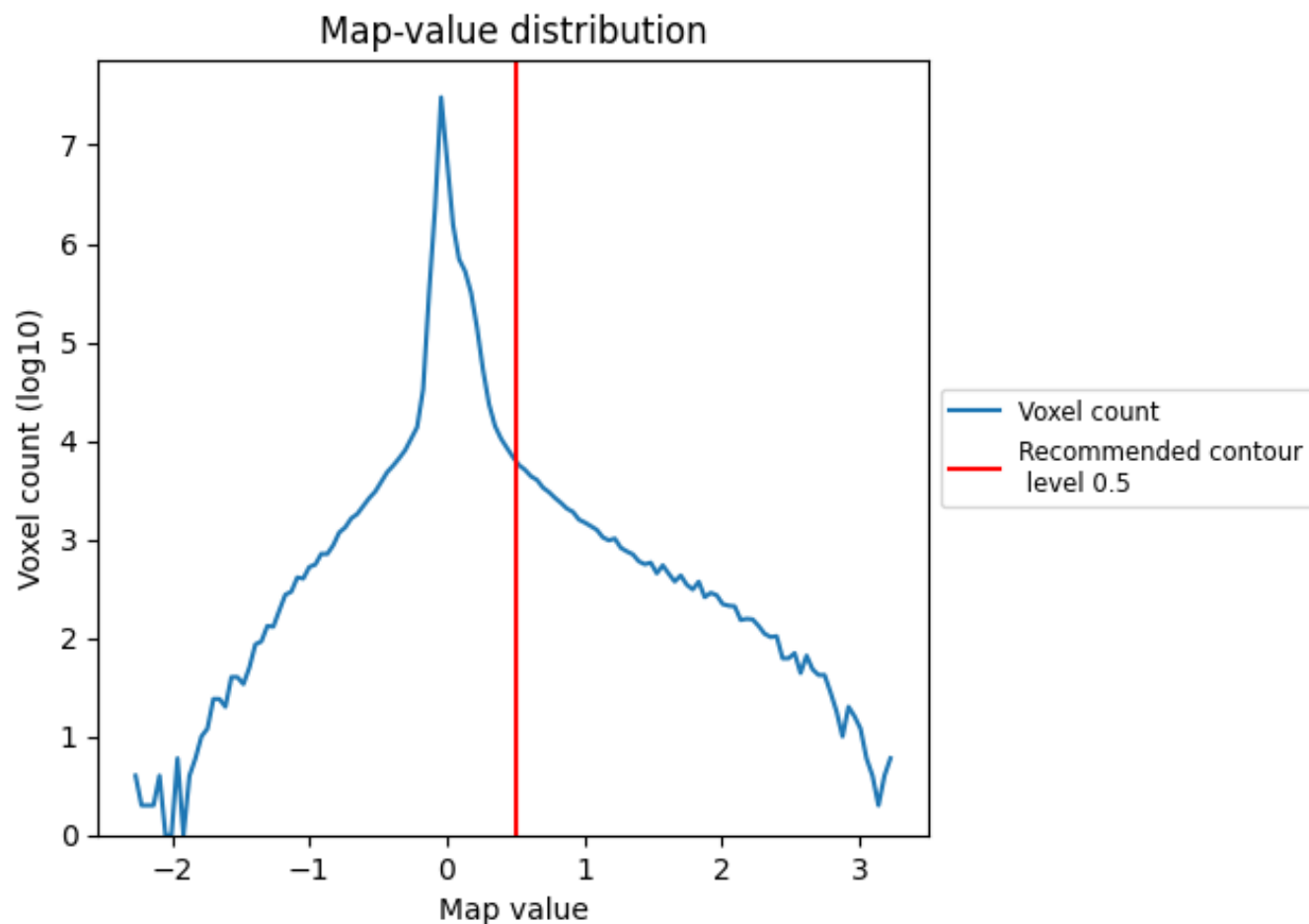
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

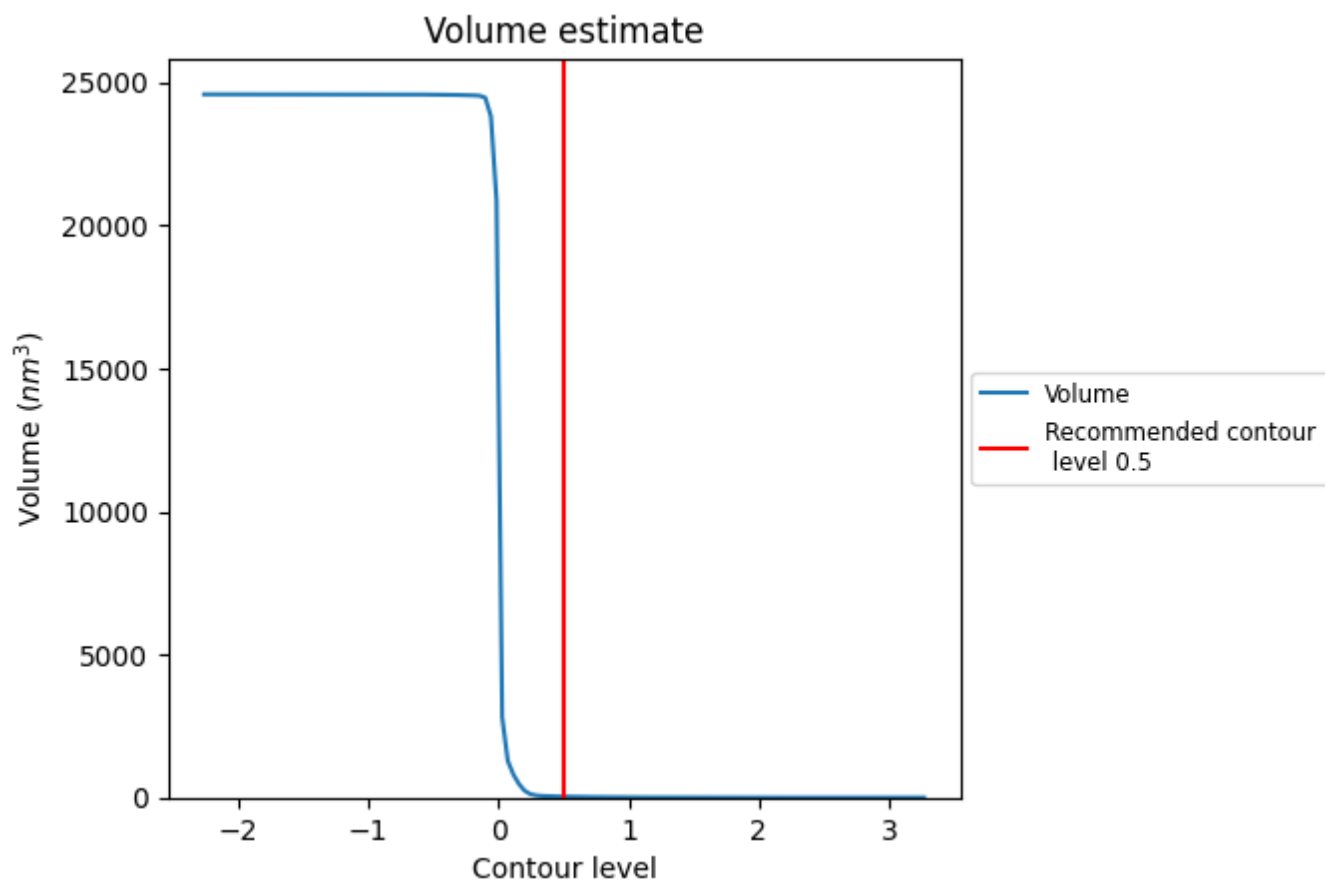
This section contains the results of statistical analysis of the map.

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.

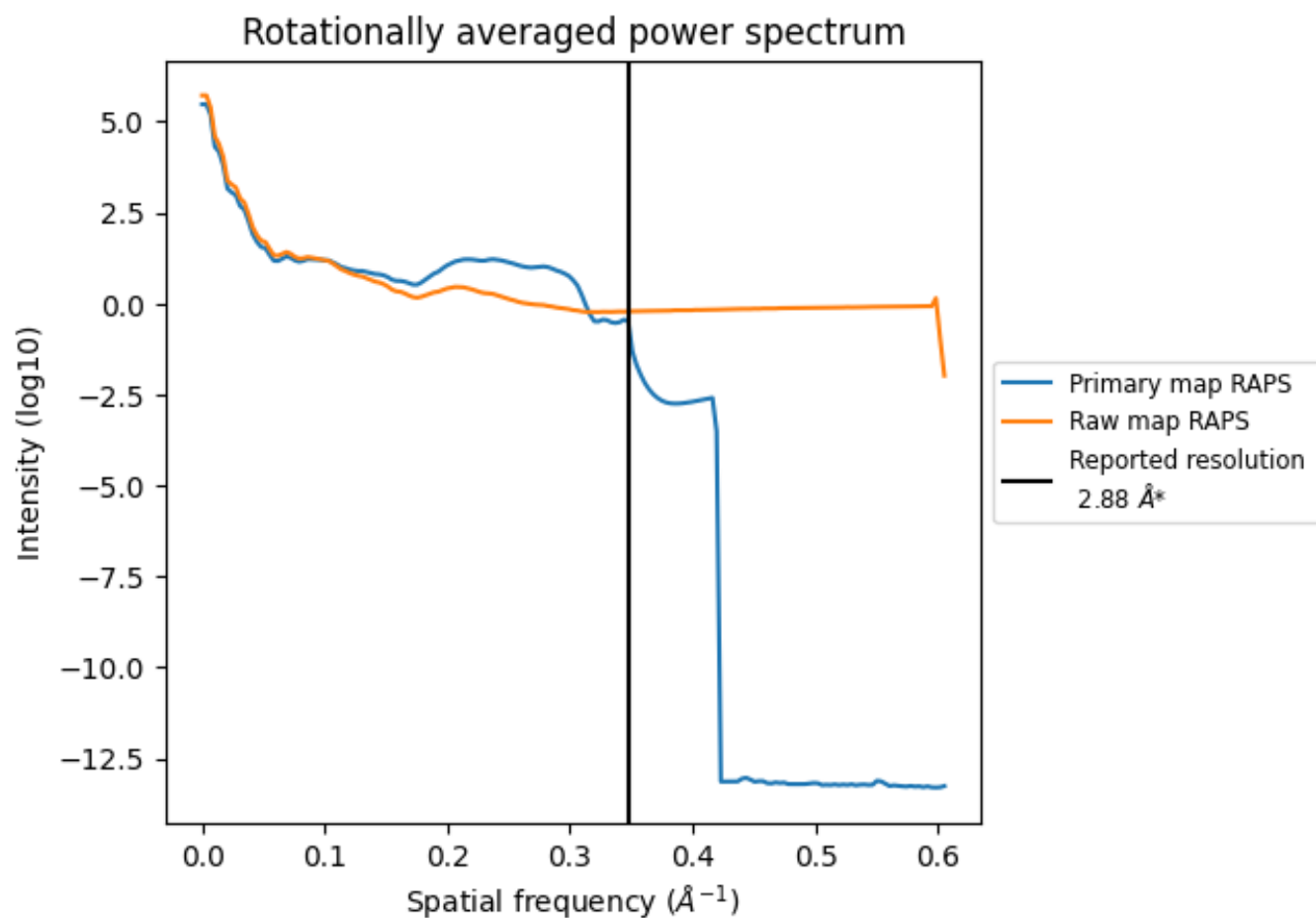
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 32 nm³; this corresponds to an approximate mass of 29 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 ⓘ

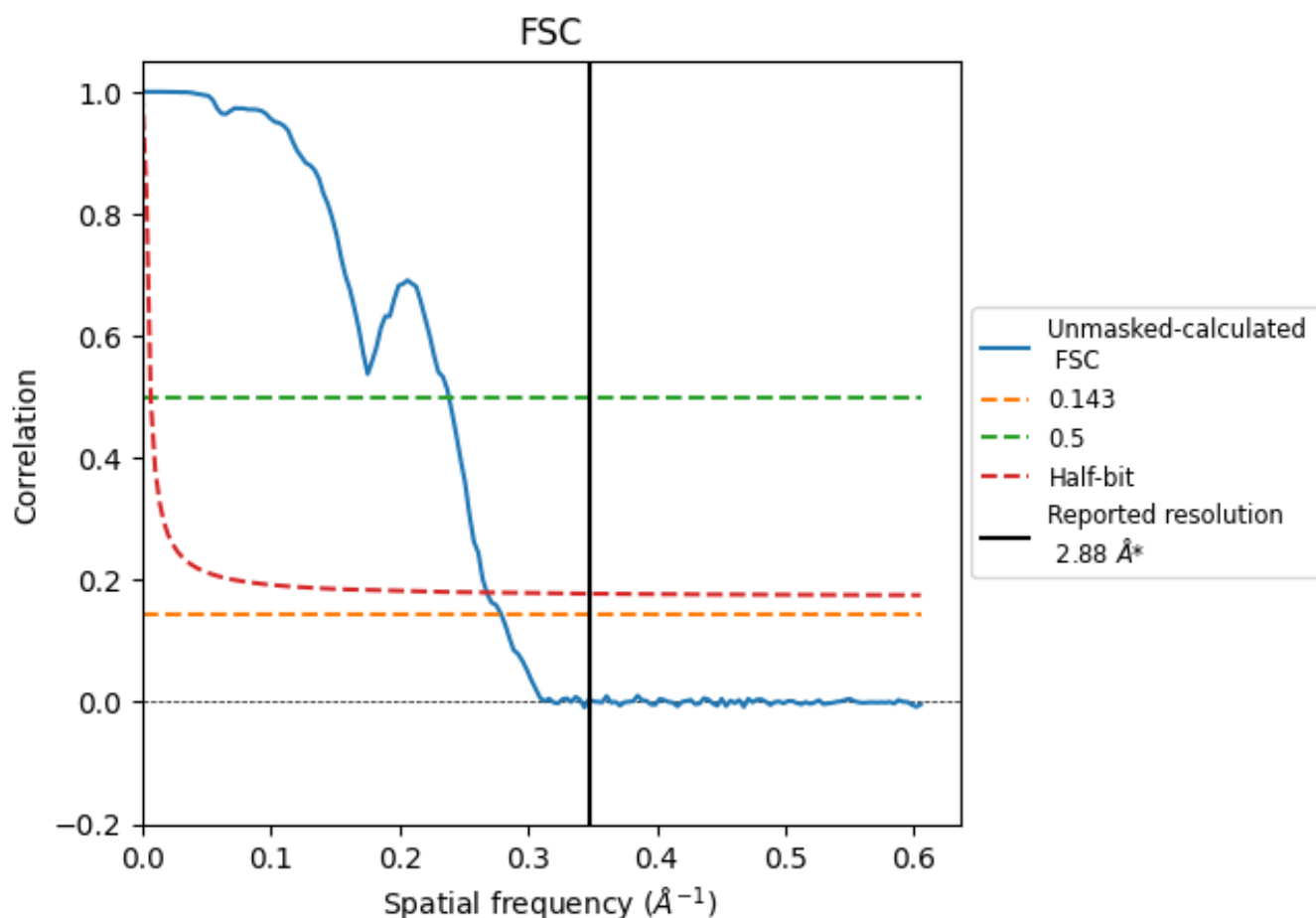


*Reported resolution corresponds to spatial frequency of 0.347 Å⁻¹

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 $0.347 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

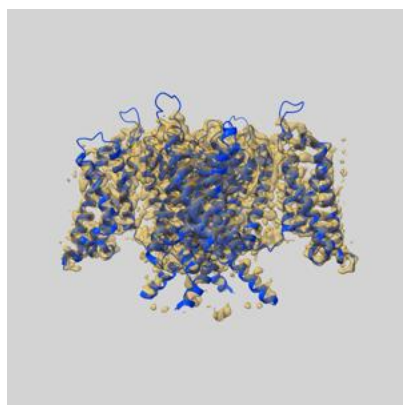
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.88	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.58	4.19	3.73

*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 3.58 differs from the reported value 2.88 by more than 10 %

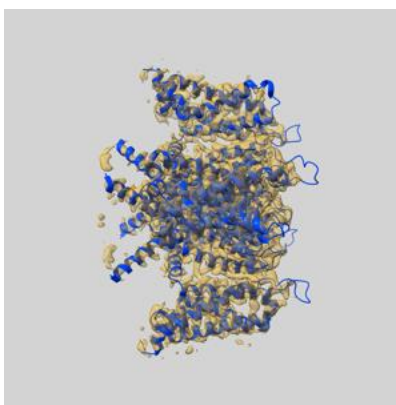
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26577 and PDB model 7UKD. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

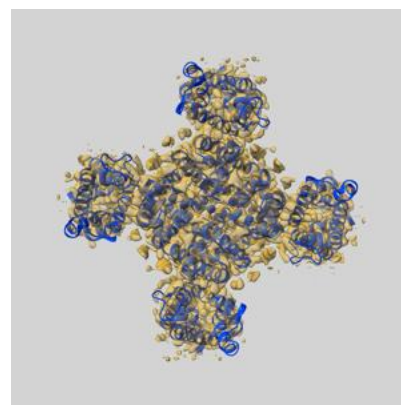
9.1 Map-model overlay [i](#)



X



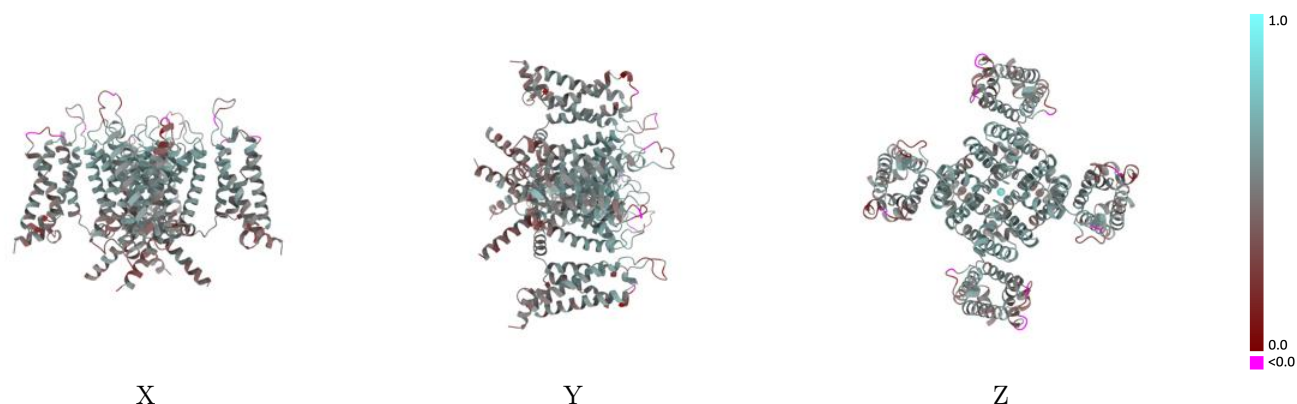
Y



Z

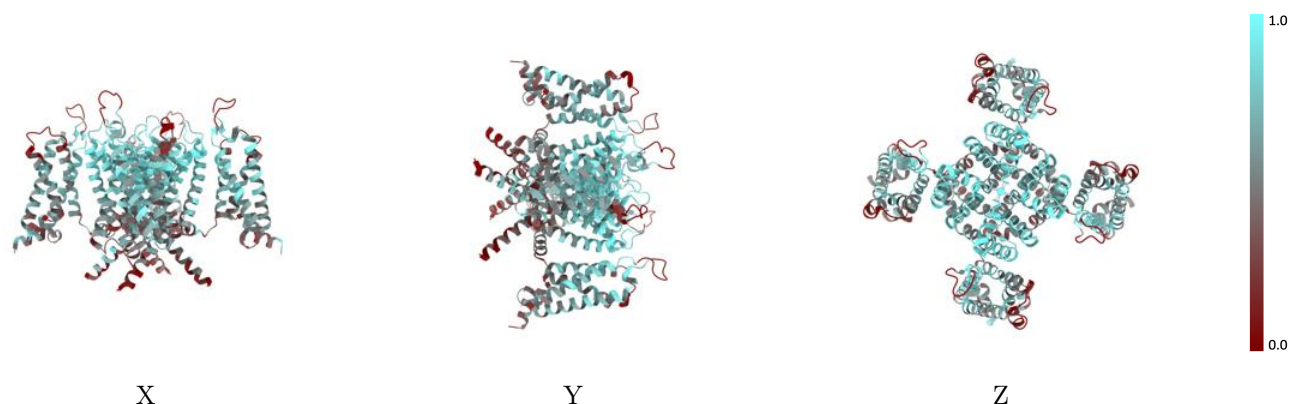
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



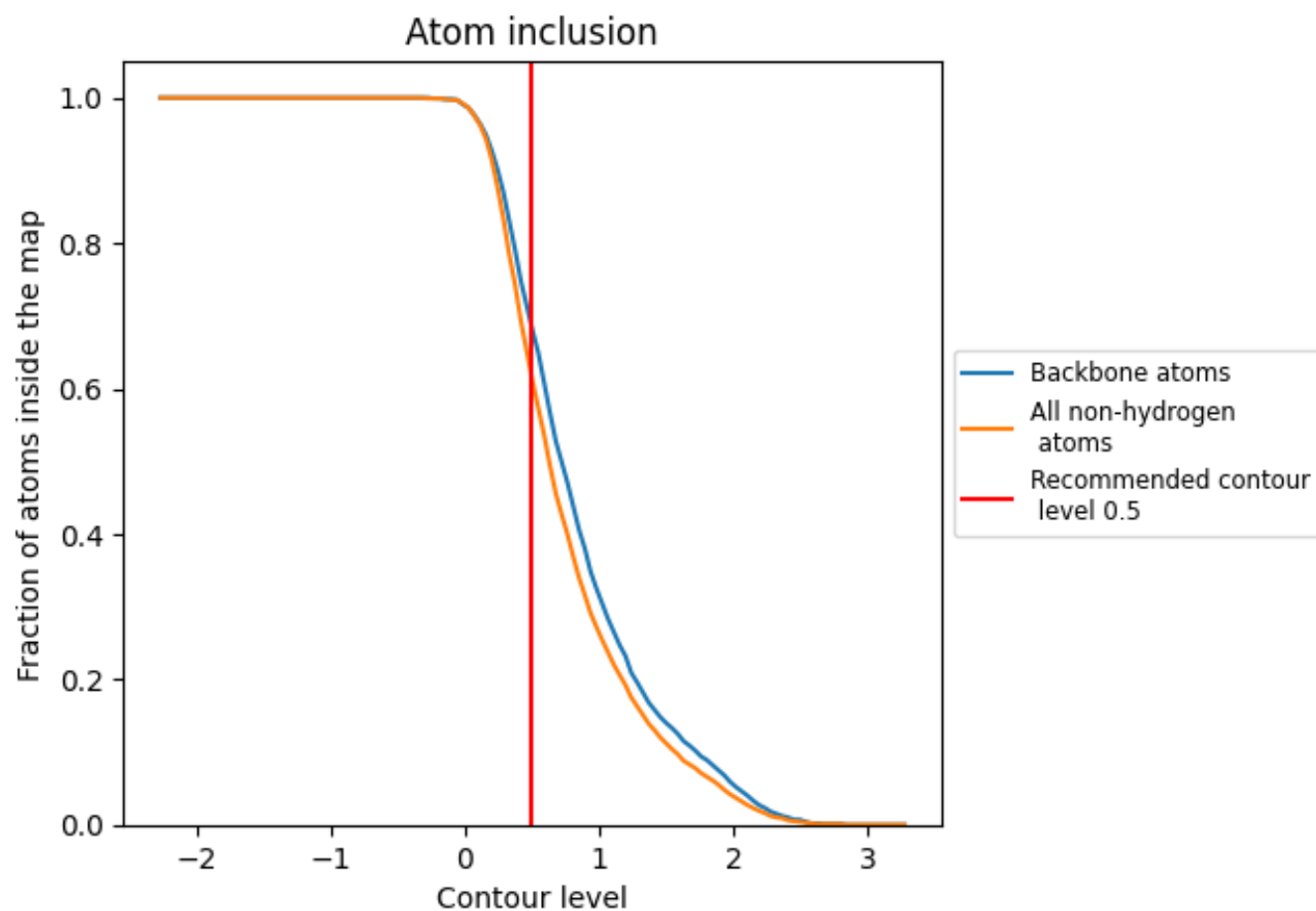
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 68% of all backbone atoms, 61% 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	0.6120	0.4940
A	0.6090	0.4910
B	0.6180	0.4960
C	0.6050	0.4900
D	0.6150	0.4970

