



Full wwPDB EM Validation Report ⓘ

Mar 25, 2023 – 12:54 PM EDT

PDB ID : 8CX2
EMDB ID : EMD-27034
Title : Cryo-EM structure of human APOBEC3G/HIV-1 Vif/CBFbeta/ELOB/ELO C dimeric complex in State 2
Authors : Li, Y.; Langley, C.; Azumaya, C.M.; Echeverria, I.; Chesarino, N.M.; Emerman, M.; Cheng, Y.; Gross, J.D.
Deposited on : 2022-05-19
Resolution : 3.20 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:

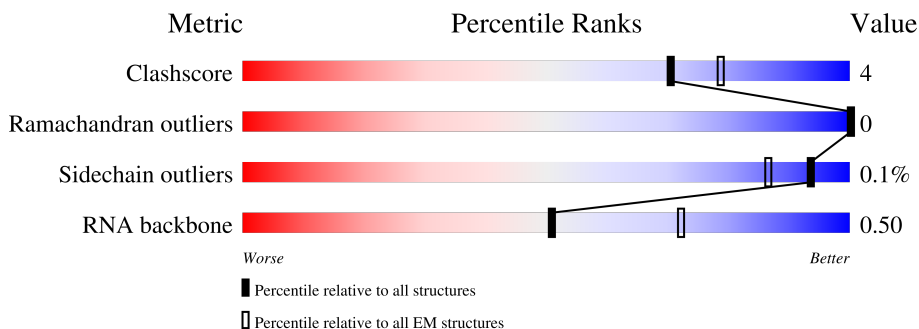
EMDB validation analysis : 0.0.1.dev50
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.32.2

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 3.20 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	421	83% 6% 10%
1	F	421	81% 8% 10%
2	B	192	85% 6% 8%
2	G	192	14% 83% 8% 8%
3	C	182	74% 8% 18%
3	H	182	6% 76% 8% 16%
4	D	118	11% 68% 7% 25%

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Mol	Chain	Length	Quality of chain
4	I	118	<p>64% 64% 10% 26%</p>
5	E	112	<p>7% 63% 13% 23%</p>
5	J	112	<p>66% 69% 8% 23%</p>
6	K	9	<p>89% 11%</p>
7	L	9	<p>78% 11% 11%</p>

2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 15041 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 DNA dC->dU-editing enzyme APOBEC-3G.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	377	3211	2064	562	560	25	0	0
1	F	377	3211	2064	562	560	25	0	0

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	385	GLY	-	expression tag	UNP Q9HC16
A	386	SER	-	expression tag	UNP Q9HC16
A	387	SER	-	expression tag	UNP Q9HC16
A	388	LEU	-	expression tag	UNP Q9HC16
A	389	GLU	-	expression tag	UNP Q9HC16
A	390	GLY	-	expression tag	UNP Q9HC16
A	391	GLY	-	expression tag	UNP Q9HC16
A	392	GLY	-	expression tag	UNP Q9HC16
A	393	GLY	-	expression tag	UNP Q9HC16
A	394	TRP	-	expression tag	UNP Q9HC16
A	395	SER	-	expression tag	UNP Q9HC16
A	396	HIS	-	expression tag	UNP Q9HC16
A	397	PRO	-	expression tag	UNP Q9HC16
A	398	GLN	-	expression tag	UNP Q9HC16
A	399	PHE	-	expression tag	UNP Q9HC16
A	400	GLU	-	expression tag	UNP Q9HC16
A	401	LYS	-	expression tag	UNP Q9HC16
A	402	GLY	-	expression tag	UNP Q9HC16
A	403	GLY	-	expression tag	UNP Q9HC16
A	404	GLY	-	expression tag	UNP Q9HC16
A	405	SER	-	expression tag	UNP Q9HC16
A	406	GLY	-	expression tag	UNP Q9HC16
A	407	GLY	-	expression tag	UNP Q9HC16
A	408	GLY	-	expression tag	UNP Q9HC16
A	409	SER	-	expression tag	UNP Q9HC16
A	410	GLY	-	expression tag	UNP Q9HC16

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Chain	Residue	Modelled	Actual	Comment	Reference
A	411	GLY	-	expression tag	UNP Q9HC16
A	412	GLY	-	expression tag	UNP Q9HC16
A	413	SER	-	expression tag	UNP Q9HC16
A	414	TRP	-	expression tag	UNP Q9HC16
A	415	SER	-	expression tag	UNP Q9HC16
A	416	HIS	-	expression tag	UNP Q9HC16
A	417	PRO	-	expression tag	UNP Q9HC16
A	418	GLN	-	expression tag	UNP Q9HC16
A	419	PHE	-	expression tag	UNP Q9HC16
A	420	GLU	-	expression tag	UNP Q9HC16
A	421	LYS	-	expression tag	UNP Q9HC16
F	385	GLY	-	expression tag	UNP Q9HC16
F	386	SER	-	expression tag	UNP Q9HC16
F	387	SER	-	expression tag	UNP Q9HC16
F	388	LEU	-	expression tag	UNP Q9HC16
F	389	GLU	-	expression tag	UNP Q9HC16
F	390	GLY	-	expression tag	UNP Q9HC16
F	391	GLY	-	expression tag	UNP Q9HC16
F	392	GLY	-	expression tag	UNP Q9HC16
F	393	GLY	-	expression tag	UNP Q9HC16
F	394	TRP	-	expression tag	UNP Q9HC16
F	395	SER	-	expression tag	UNP Q9HC16
F	396	HIS	-	expression tag	UNP Q9HC16
F	397	PRO	-	expression tag	UNP Q9HC16
F	398	GLN	-	expression tag	UNP Q9HC16
F	399	PHE	-	expression tag	UNP Q9HC16
F	400	GLU	-	expression tag	UNP Q9HC16
F	401	LYS	-	expression tag	UNP Q9HC16
F	402	GLY	-	expression tag	UNP Q9HC16
F	403	GLY	-	expression tag	UNP Q9HC16
F	404	GLY	-	expression tag	UNP Q9HC16
F	405	SER	-	expression tag	UNP Q9HC16
F	406	GLY	-	expression tag	UNP Q9HC16
F	407	GLY	-	expression tag	UNP Q9HC16
F	408	GLY	-	expression tag	UNP Q9HC16
F	409	SER	-	expression tag	UNP Q9HC16
F	410	GLY	-	expression tag	UNP Q9HC16
F	411	GLY	-	expression tag	UNP Q9HC16
F	412	GLY	-	expression tag	UNP Q9HC16
F	413	SER	-	expression tag	UNP Q9HC16
F	414	TRP	-	expression tag	UNP Q9HC16
F	415	SER	-	expression tag	UNP Q9HC16

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Chain	Residue	Modelled	Actual	Comment	Reference
F	416	HIS	-	expression tag	UNP Q9HC16
F	417	PRO	-	expression tag	UNP Q9HC16
F	418	GLN	-	expression tag	UNP Q9HC16
F	419	PHE	-	expression tag	UNP Q9HC16
F	420	GLU	-	expression tag	UNP Q9HC16
F	421	LYS	-	expression tag	UNP Q9HC16

- Molecule 2 is a protein called Virion infectivity factor.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	176	1468	945	270	247	6	0	0
2	G	176	1468	945	270	247	6	0	0

- Molecule 3 is a protein called Core-binding factor subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	150	1254	776	234	238	6	0	0
3	H	153	1275	791	237	241	6	0	0

- Molecule 4 is a protein called Elongin-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	88	698	442	121	132	3	0	0
4	I	87	690	436	120	131	3	0	0

- Molecule 5 is a protein called Elongin-C.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	86	688	444	110	128	6	0	0
5	J	86	688	444	110	128	6	0	0

- Molecule 6 is a RNA chain called RNA (5'-R(P*AP*AP*UP*AP*AP*AP*AP*U)-3').

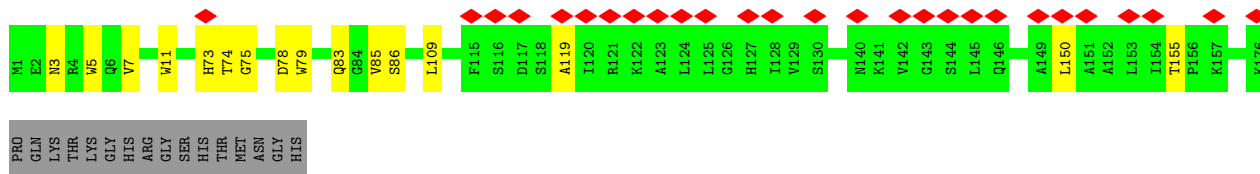
Mol	Chain	Residues	Atoms				AltConf	Trace	
			Total	C	N	O			P
6	K	9	194	88	39	58	9	0	0

- Molecule 7 is a RNA chain called RNA (5'-R(P*UP*UP*UP*AP*AP*AP*AP*U)-3').

Mol	Chain	Residues	Atoms				AltConf	Trace	
			Total	C	N	O			P
7	L	9	190	86	33	62	9	0	0

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

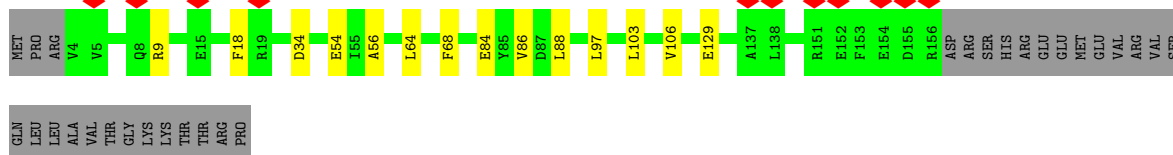
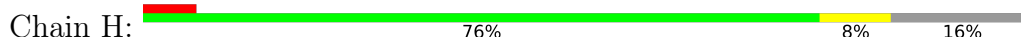
Mol	Chain	Residues	Atoms		AltConf
8	A	2	Total	Zn	0
			2	2	
8	B	1	Total	Zn	0
			1	1	
8	F	2	Total	Zn	0
			2	2	
8	G	1	Total	Zn	0
			1	1	



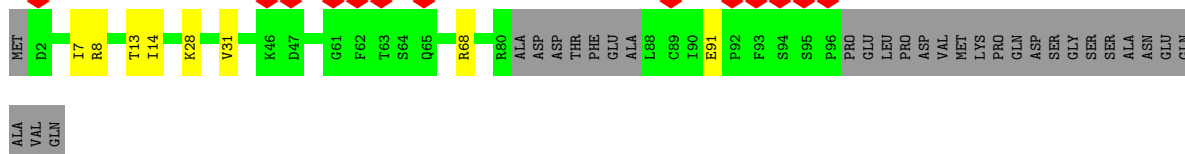
● Molecule 3: Core-binding factor subunit beta



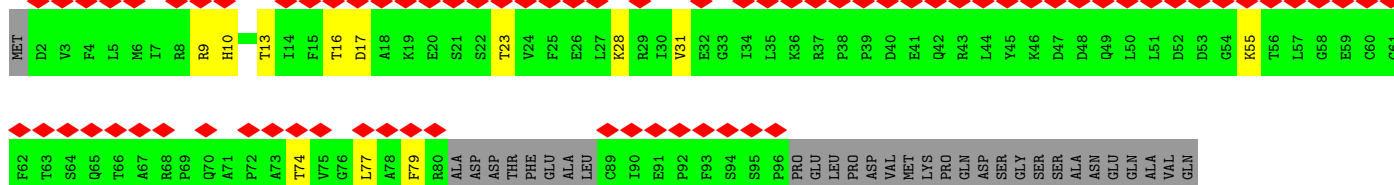
● Molecule 3: Core-binding factor subunit beta



● Molecule 4: Elongin-B

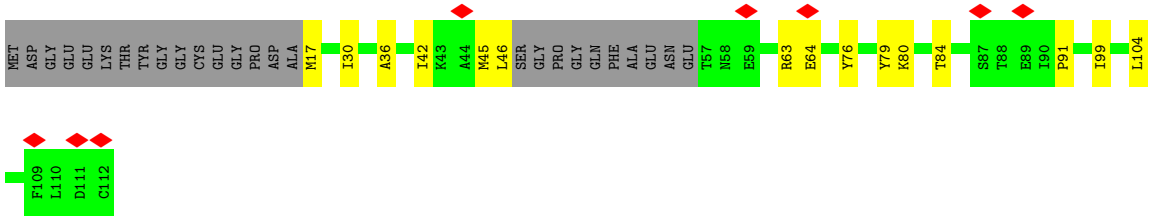


● Molecule 4: Elongin-B

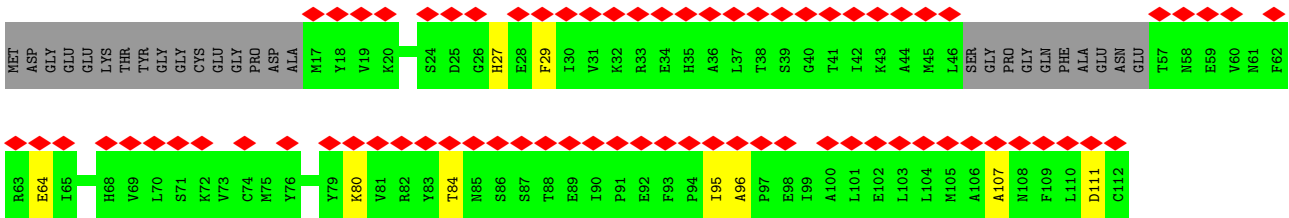


● Molecule 5: Elongin-C

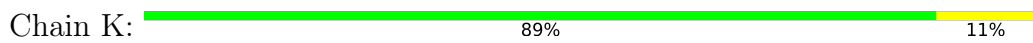




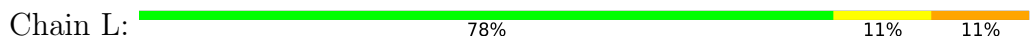
• Molecule 5: Elongin-C



• Molecule 6: RNA (5'-R(P*AP*AP*UP*AP*AP*AP*AP*AP*U)-3')



• Molecule 7: RNA (5'-R(P*UP*UP*UP*AP*AP*AP*AP*AP*U)-3')



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	46083	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 ($e^-/\text{\AA}^2$)	68	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.401	Depositor
Minimum map value	-0.621	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.039	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	293.91998, 293.91998, 293.91998	wwPDB
Map dimensions	352, 352, 352	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.835, 0.835, 0.835	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:
ZN

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	0.28	0/3318	0.50	0/4496
1	F	0.27	0/3318	0.50	0/4496
2	B	0.32	0/1515	0.54	0/2055
2	G	0.32	0/1515	0.52	0/2055
3	C	0.26	0/1278	0.57	0/1712
3	H	0.25	0/1300	0.56	0/1744
4	D	0.25	0/710	0.52	0/956
4	I	0.25	0/702	0.51	0/945
5	E	0.24	0/702	0.44	0/947
5	J	0.27	0/702	0.46	0/947
6	K	0.21	0/218	0.65	0/337
7	L	0.12	0/212	0.67	0/327
All	All	0.28	0/15490	0.52	0/21017

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	3211	0	3057	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3211	0	3057	20	0
2	B	1468	0	1462	11	0
2	G	1468	0	1462	13	0
3	C	1254	0	1191	14	0
3	H	1275	0	1216	18	0
4	D	698	0	705	4	0
4	I	690	0	694	7	0
5	E	688	0	690	11	0
5	J	688	0	690	6	0
6	K	194	0	98	2	0
7	L	190	0	96	2	0
8	A	2	0	0	0	0
8	B	1	0	0	0	0
8	F	2	0	0	0	0
8	G	1	0	0	0	0
All	All	15041	0	14418	107	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (107) 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)
2:B:145:LEU:HD21	5:E:76:TYR:CG	2.24	0.72
5:E:64:GLU:N	5:E:64:GLU:OE1	2.27	0.67
1:F:259:GLU:N	1:F:259:GLU:OE1	2.33	0.62
1:A:259:GLU:N	1:A:259:GLU:OE1	2.35	0.57
2:G:73:HIS:HB3	3:H:56:ALA:HB2	1.87	0.57
1:F:120:VAL:HG22	1:F:122:ARG:O	2.05	0.55
1:F:227:MET:HE2	1:F:232:TRP:CE3	2.41	0.55
1:F:327:THR:HA	1:F:330:GLU:HG2	1.87	0.55
5:E:17:MET:SD	5:E:17:MET:N	2.81	0.54
1:F:321:CYS:O	1:F:325:LEU:HD13	2.08	0.54
1:A:337:ILE:HD11	1:A:375:LEU:HD22	1.90	0.53
2:B:109:LEU:HA	3:C:151:ARG:HB2	1.91	0.53
3:H:86:VAL:HG13	3:H:97:LEU:CD2	2.38	0.53
4:D:8:ARG:NH2	4:D:91:GLU:O	2.41	0.53
3:H:86:VAL:HG22	3:H:97:LEU:CD2	2.39	0.52
3:H:18:PHE:CE2	3:H:103:LEU:HD23	2.44	0.52
4:I:13:THR:OG1	5:J:27:HIS:NE2	2.39	0.52
1:F:14:ARG:NH1	1:F:170:GLU:OE2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:8:GLN:OE1	3:C:106:VAL:HG23	2.09	0.52
1:A:119:PHE:HB3	1:A:152:MET:HG2	1.92	0.51
1:F:191:GLU:OE2	1:F:222:TYR:OH	2.27	0.50
5:E:63:ARG:NE	5:E:63:ARG:O	2.45	0.50
2:G:119:ALA:HB1	2:G:150:LEU:HD21	1.92	0.50
3:C:54:GLU:N	3:C:54:GLU:OE1	2.42	0.50
2:G:7:VAL:HG22	3:H:68:PHE:CD2	2.45	0.50
1:A:14:ARG:NH1	1:A:168:GLN:OE1	2.45	0.50
5:E:79:TYR:OH	5:E:91:PRO:O	2.26	0.50
1:A:59:TYR:HB3	6:K:3:A:C8	2.47	0.49
1:A:180:LYS:O	1:A:183:ILE:HG22	2.13	0.48
4:I:16:THR:HG22	4:I:17:ASP:N	2.28	0.48
4:I:28:LYS:HA	4:I:31:VAL:HG12	1.95	0.48
1:F:19:TYR:O	1:F:30:ARG:NE	2.46	0.48
4:I:10:HIS:O	4:I:10:HIS:ND1	2.47	0.48
3:H:84:GLU:OE1	3:H:84:GLU:N	2.42	0.48
3:H:86:VAL:HG22	3:H:97:LEU:HD23	1.96	0.47
1:A:187:ILE:HG21	1:A:265:VAL:HG21	1.97	0.47
3:H:103:LEU:O	3:H:106:VAL:HG12	2.13	0.47
1:A:80:LEU:HD21	2:G:155:THR:HG23	1.97	0.47
2:B:81:LEU:HD21	2:B:174:TRP:CZ3	2.49	0.47
1:F:262:PHE:O	1:F:266:ILE:HG12	2.14	0.47
5:J:64:GLU:N	5:J:64:GLU:OE1	2.47	0.46
1:A:225:GLU:OE1	1:A:278:ARG:NH2	2.48	0.46
1:A:100:CYS:O	1:A:104:MET:HG3	2.15	0.46
1:A:187:ILE:HD13	1:A:265:VAL:HG21	1.97	0.46
4:I:23:THR:OG1	4:I:55:LYS:O	2.30	0.46
3:C:86:VAL:HG13	3:C:97:LEU:HD23	1.96	0.46
1:A:337:ILE:CD1	1:A:375:LEU:HD22	2.44	0.46
5:E:99:ILE:O	5:E:99:ILE:HG22	2.16	0.46
5:J:107:ALA:O	5:J:111:ASP:N	2.49	0.46
2:G:78:ASP:O	2:G:79:TRP:HB2	2.17	0.45
3:H:9:ARG:NH2	3:H:129:GLU:OE2	2.48	0.45
3:C:86:VAL:HG13	3:C:97:LEU:CD2	2.46	0.45
2:G:109:LEU:HD23	2:G:109:LEU:O	2.16	0.45
1:F:262:PHE:O	1:F:265:VAL:HG12	2.17	0.45
2:G:73:HIS:HB2	3:H:54:GLU:HG2	1.99	0.45
1:F:262:PHE:HZ	1:F:279:VAL:HG11	1.81	0.45
2:B:9:ILE:HD13	3:C:18:PHE:CE2	2.52	0.44
2:B:77:ARG:NE	3:C:34:ASP:HA	2.32	0.44
1:F:132:GLN:HG2	1:F:192:ILE:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:86:VAL:HG13	3:H:97:LEU:HD23	1.98	0.44
3:C:108:VAL:HG21	3:C:125:LEU:HD11	2.00	0.44
5:E:45:MET:HB3	5:E:46:LEU:HD12	2.00	0.44
2:G:3:ASN:CB	2:G:5:TRP:HE1	2.31	0.44
1:A:311:THR:HG22	1:A:312:ALA:N	2.32	0.44
1:A:55:ARG:NH1	7:L:6:U:O4'	2.52	0.43
4:I:74:THR:O	4:I:74:THR:HG23	2.18	0.43
5:J:80:LYS:O	5:J:84:THR:N	2.51	0.43
5:J:27:HIS:HD2	5:J:29:PHE:CE1	2.36	0.43
1:A:227:MET:HE3	1:A:227:MET:HA	2.01	0.43
4:D:28:LYS:HA	4:D:31:VAL:HG12	1.99	0.43
3:H:88:LEU:HD23	3:H:88:LEU:O	2.19	0.43
3:H:34:ASP:C	3:H:34:ASP:OD2	2.57	0.43
1:A:262:PHE:O	1:A:266:ILE:HG13	2.19	0.42
2:B:147:TYR:HB2	3:C:152:GLU:OE1	2.19	0.42
5:E:36:ALA:HB1	5:E:42:ILE:HG21	2.01	0.42
1:F:207:ASN:HA	1:F:361:TRP:CZ2	2.54	0.42
2:B:149:ALA:HB1	5:E:104:LEU:HB2	2.00	0.42
2:B:20:THR:HG23	6:K:3:A:H61	1.84	0.42
3:H:54:GLU:OE1	3:H:54:GLU:N	2.43	0.42
1:F:24:ARG:NH2	7:L:0:U:OP2	2.49	0.42
3:C:18:PHE:CD2	3:C:103:LEU:HD23	2.54	0.42
3:C:19:ARG:O	3:C:19:ARG:HG2	2.20	0.42
5:E:80:LYS:O	5:E:84:THR:N	2.53	0.42
2:G:74:THR:O	2:G:75:GLY:C	2.55	0.42
2:B:105:GLN:O	2:B:109:LEU:HD13	2.20	0.42
2:B:147:TYR:CD1	3:C:152:GLU:OE2	2.73	0.42
4:D:7:ILE:O	4:D:13:THR:HA	2.20	0.42
5:J:95:ILE:HG22	5:J:96:ALA:N	2.35	0.41
1:F:180:LYS:O	1:F:183:ILE:HG22	2.19	0.41
1:F:130:ASP:OD2	2:G:83:GLN:NE2	2.49	0.41
1:F:310:PHE:HB3	1:F:338:MET:SD	2.60	0.41
3:H:54:GLU:HA	3:H:68:PHE:O	2.20	0.41
2:G:85:VAL:HG12	2:G:86:SER:N	2.35	0.41
3:H:18:PHE:CD2	3:H:103:LEU:HD23	2.56	0.41
1:A:69:ARG:HH11	1:A:69:ARG:HG2	1.86	0.41
2:G:5:TRP:CZ3	3:H:68:PHE:HB3	2.55	0.41
1:A:259:GLU:O	1:A:263:LEU:HG	2.21	0.41
1:F:199:PRO:N	1:F:200:PRO:CD	2.84	0.41
2:B:85:VAL:HG12	2:B:86:SER:N	2.36	0.41
3:C:108:VAL:CG2	3:C:125:LEU:HD11	2.51	0.40

Continued on next page...

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:11:TRP:CZ3	3:H:64:LEU:HD22	2.57	0.40
4:I:77:LEU:HD21	4:I:79:PHE:CE1	2.55	0.40
1:F:143:ASP:OD1	1:F:143:ASP:N	2.54	0.40
1:F:239:ARG:NH2	1:F:349:THR:O	2.54	0.40
1:A:104:MET:O	1:A:108:LEU:HD13	2.21	0.40
3:C:115:ASP:OD2	3:C:118:ARG:N	2.48	0.40
4:D:14:ILE:HG23	5:E:30:ILE: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	375/421 (89%)	368 (98%)	7 (2%)	0	100	100
1	F	375/421 (89%)	365 (97%)	10 (3%)	0	100	100
2	B	174/192 (91%)	167 (96%)	7 (4%)	0	100	100
2	G	174/192 (91%)	169 (97%)	5 (3%)	0	100	100
3	C	148/182 (81%)	144 (97%)	4 (3%)	0	100	100
3	H	151/182 (83%)	147 (97%)	4 (3%)	0	100	100
4	D	84/118 (71%)	83 (99%)	1 (1%)	0	100	100
4	I	83/118 (70%)	83 (100%)	0	0	100	100
5	E	82/112 (73%)	82 (100%)	0	0	100	100
5	J	82/112 (73%)	82 (100%)	0	0	100	100
All	All	1728/2050 (84%)	1690 (98%)	38 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	348/378 (92%)	348 (100%)	0	100	100
1	F	348/378 (92%)	348 (100%)	0	100	100
2	B	157/170 (92%)	157 (100%)	0	100	100
2	G	157/170 (92%)	157 (100%)	0	100	100
3	C	129/159 (81%)	129 (100%)	0	100	100
3	H	132/159 (83%)	132 (100%)	0	100	100
4	D	78/103 (76%)	77 (99%)	1 (1%)	69	87
4	I	77/103 (75%)	76 (99%)	1 (1%)	69	87
5	E	78/96 (81%)	78 (100%)	0	100	100
5	J	78/96 (81%)	78 (100%)	0	100	100
All	All	1582/1812 (87%)	1580 (100%)	2 (0%)	93	98

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

Mol	Chain	Res	Type
4	D	68	ARG
4	I	9	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
6	K	8/9 (88%)	0	0
7	L	8/9 (88%)	1 (12%)	0
All	All	16/18 (88%)	1 (6%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
7	L	6	U

There are no RNA pucker outliers to report.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

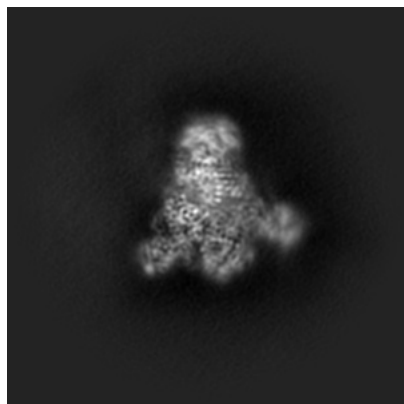
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27034. These allow visual inspection of the internal detail of the map and identification of artifacts.

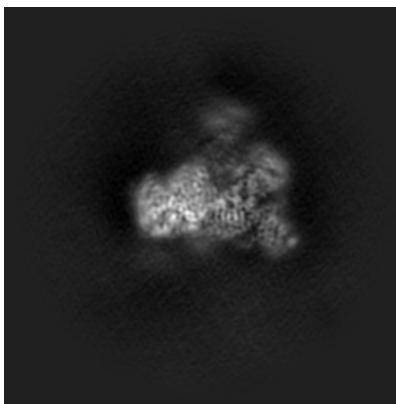
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

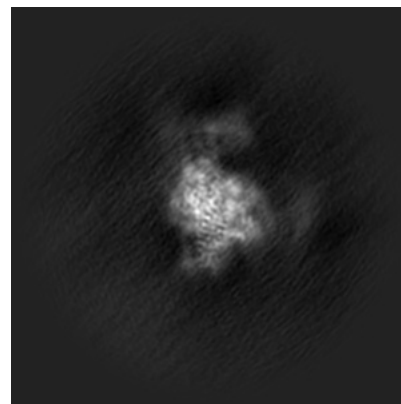
6.1.1 Primary map



X

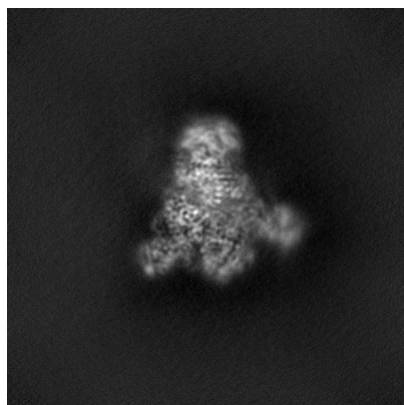


Y

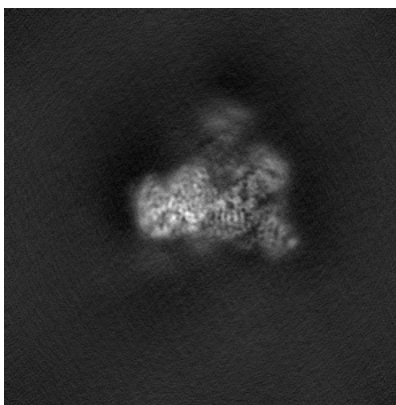


Z

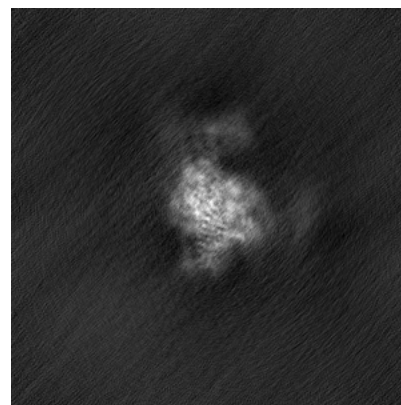
6.1.2 Raw map



X



Y

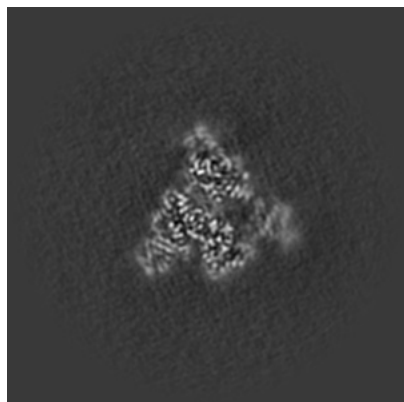


Z

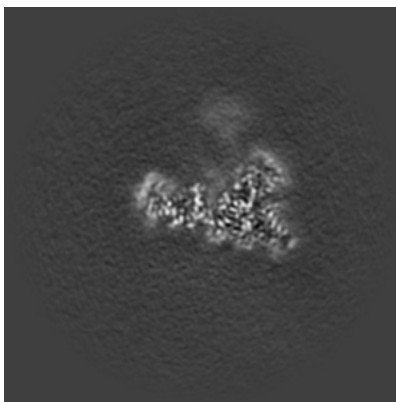
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

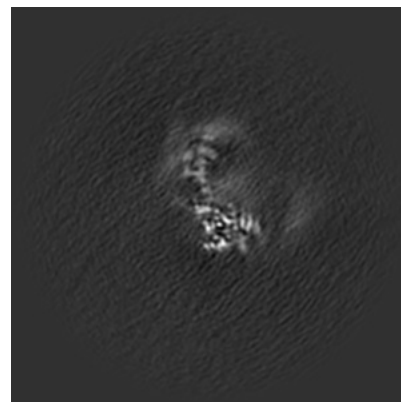
6.2.1 Primary map



X Index: 176

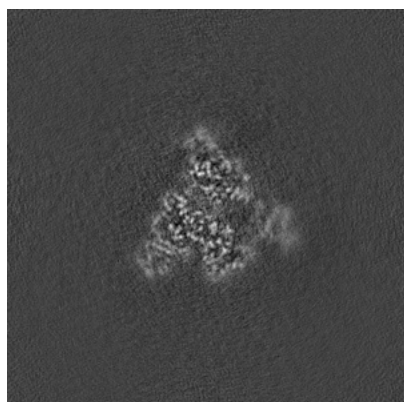


Y Index: 176

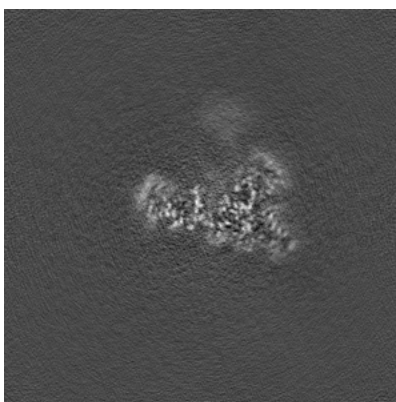


Z Index: 176

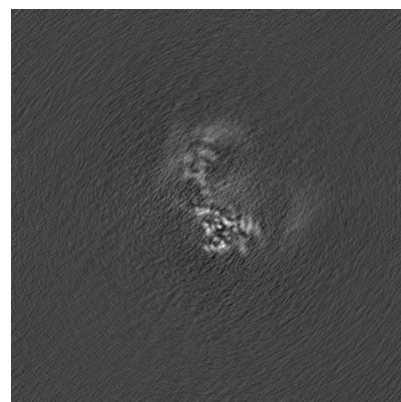
6.2.2 Raw map



X Index: 176



Y Index: 176

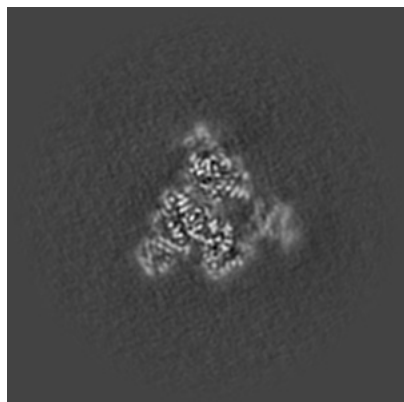


Z Index: 176

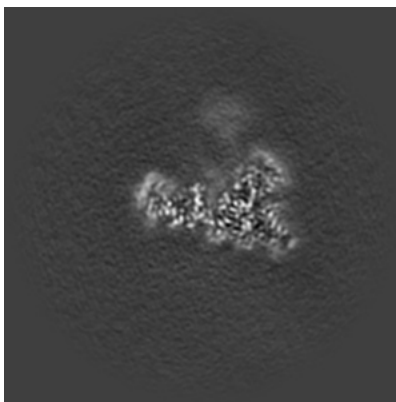
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

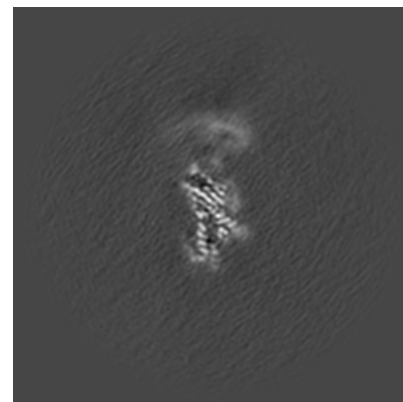
6.3.1 Primary map



X Index: 177

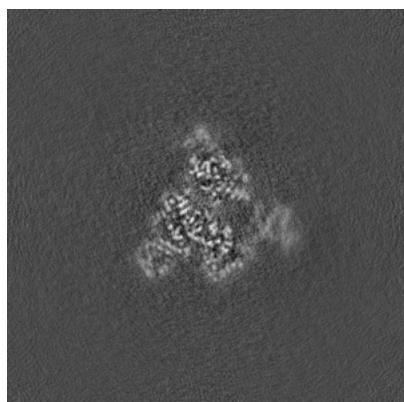


Y Index: 176

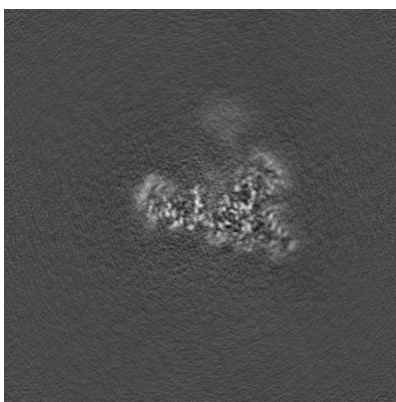


Z Index: 147

6.3.2 Raw map



X Index: 177



Y Index: 176

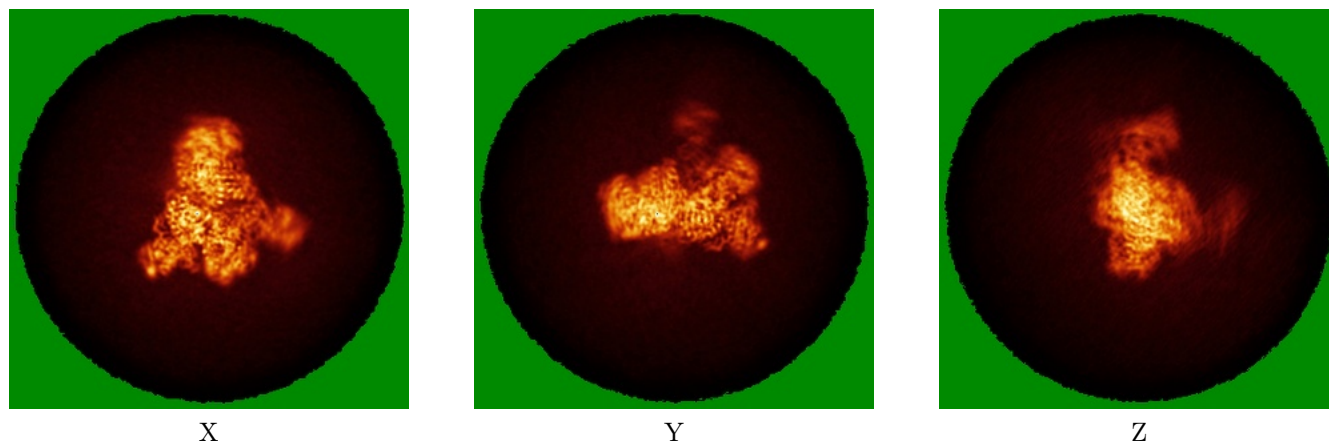


Z Index: 168

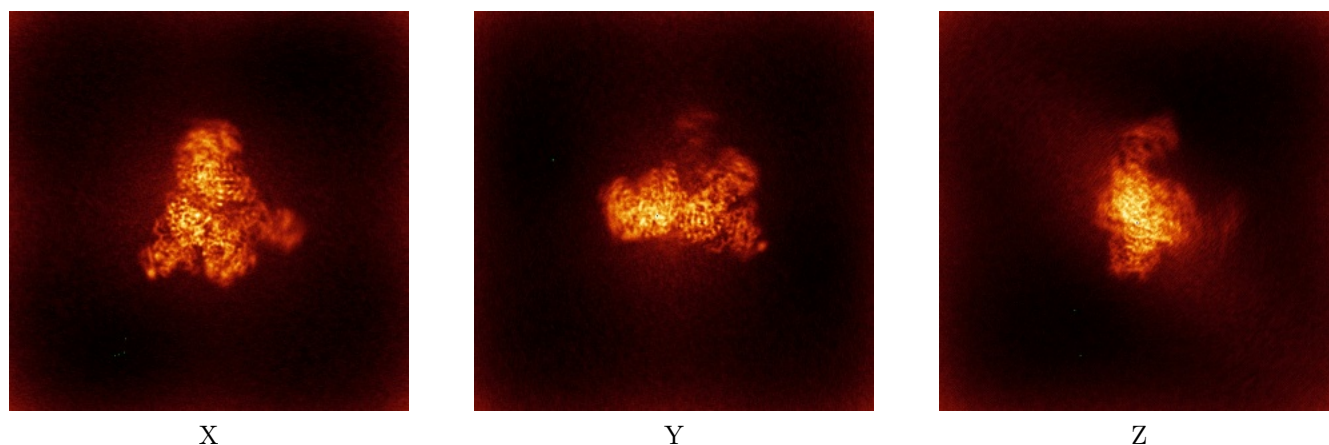
The images above show the largest variance slices of the map in three orthogonal directions.

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.

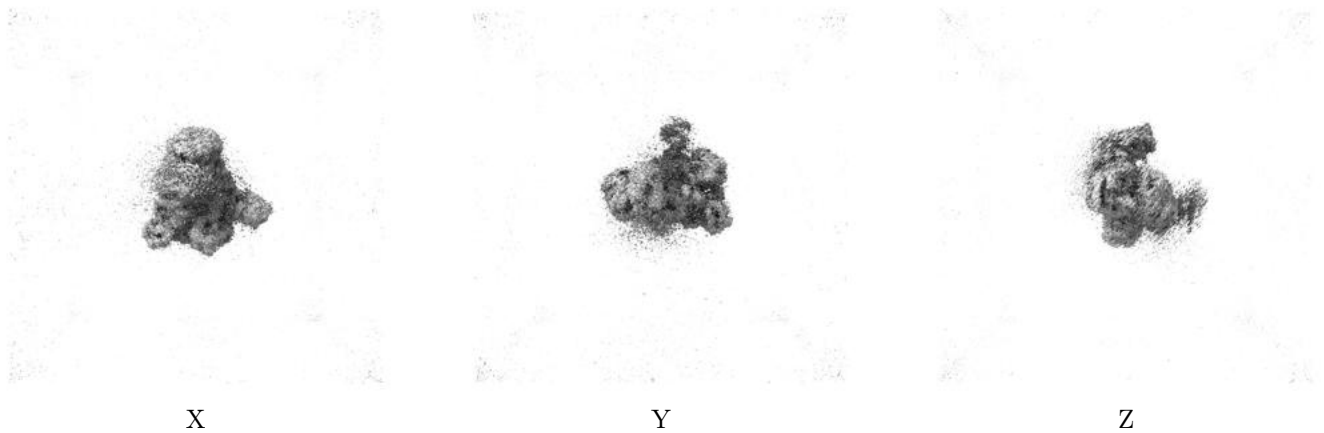
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. 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.

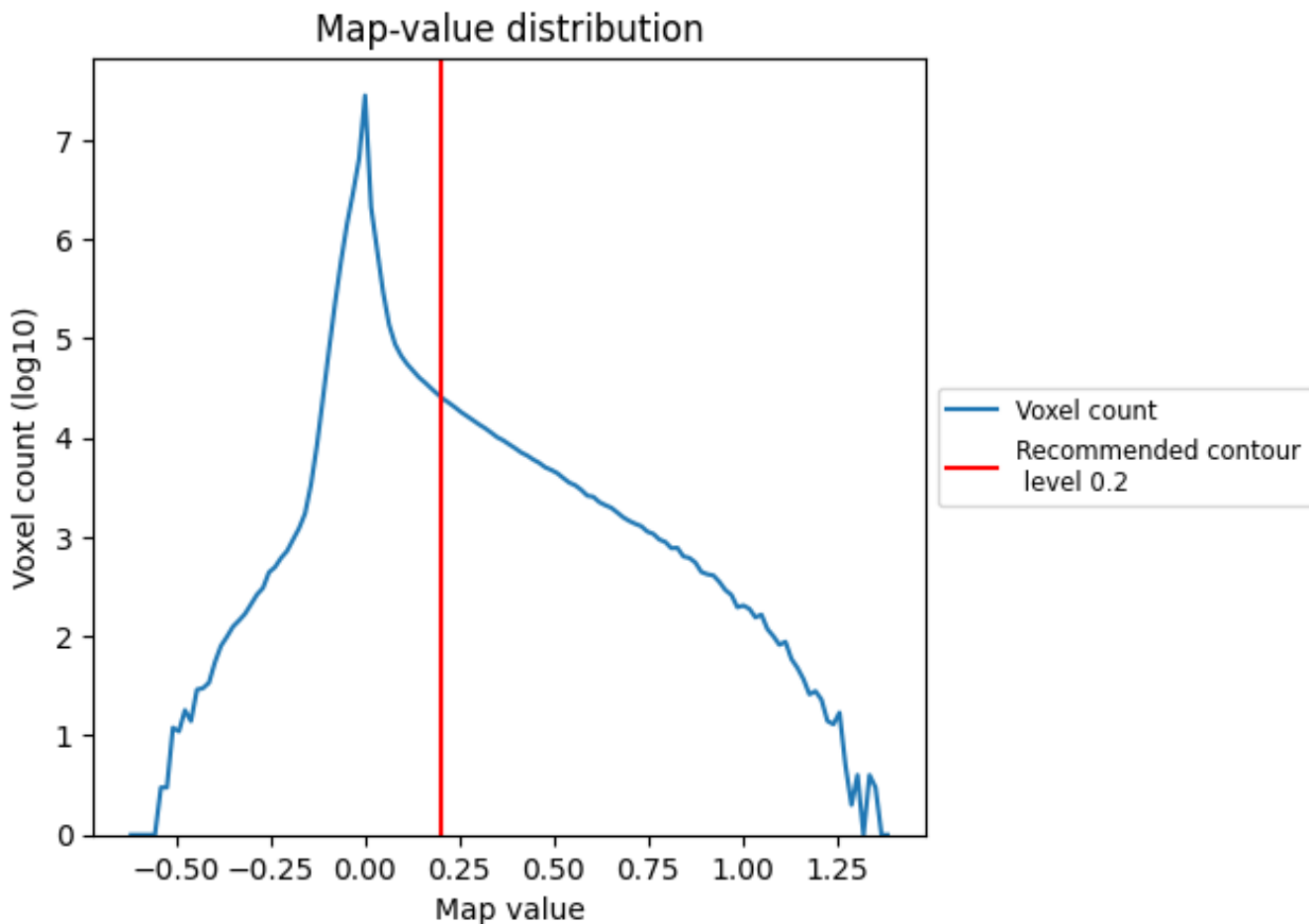
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

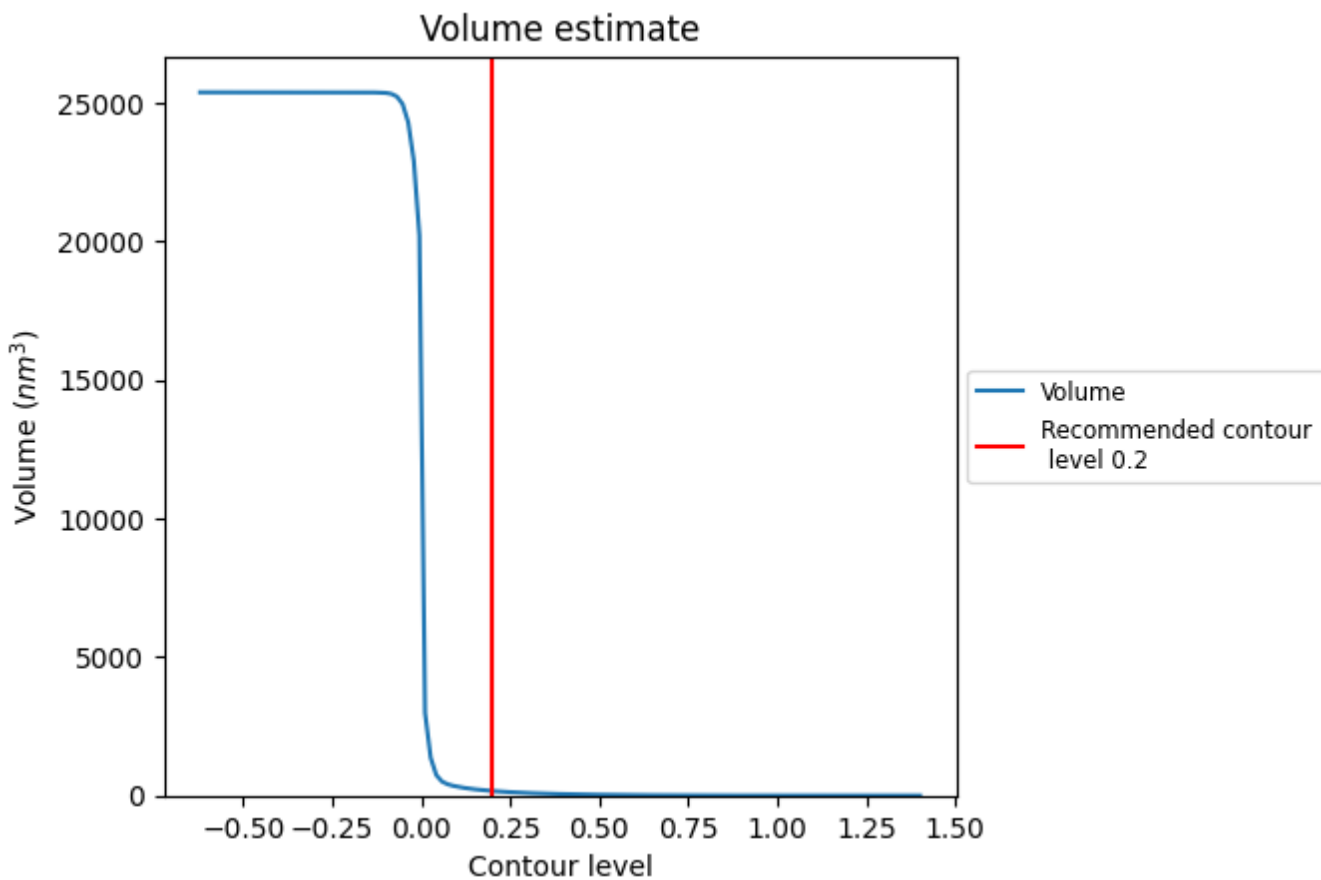
This section contains the results of statistical analysis of the map.

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.

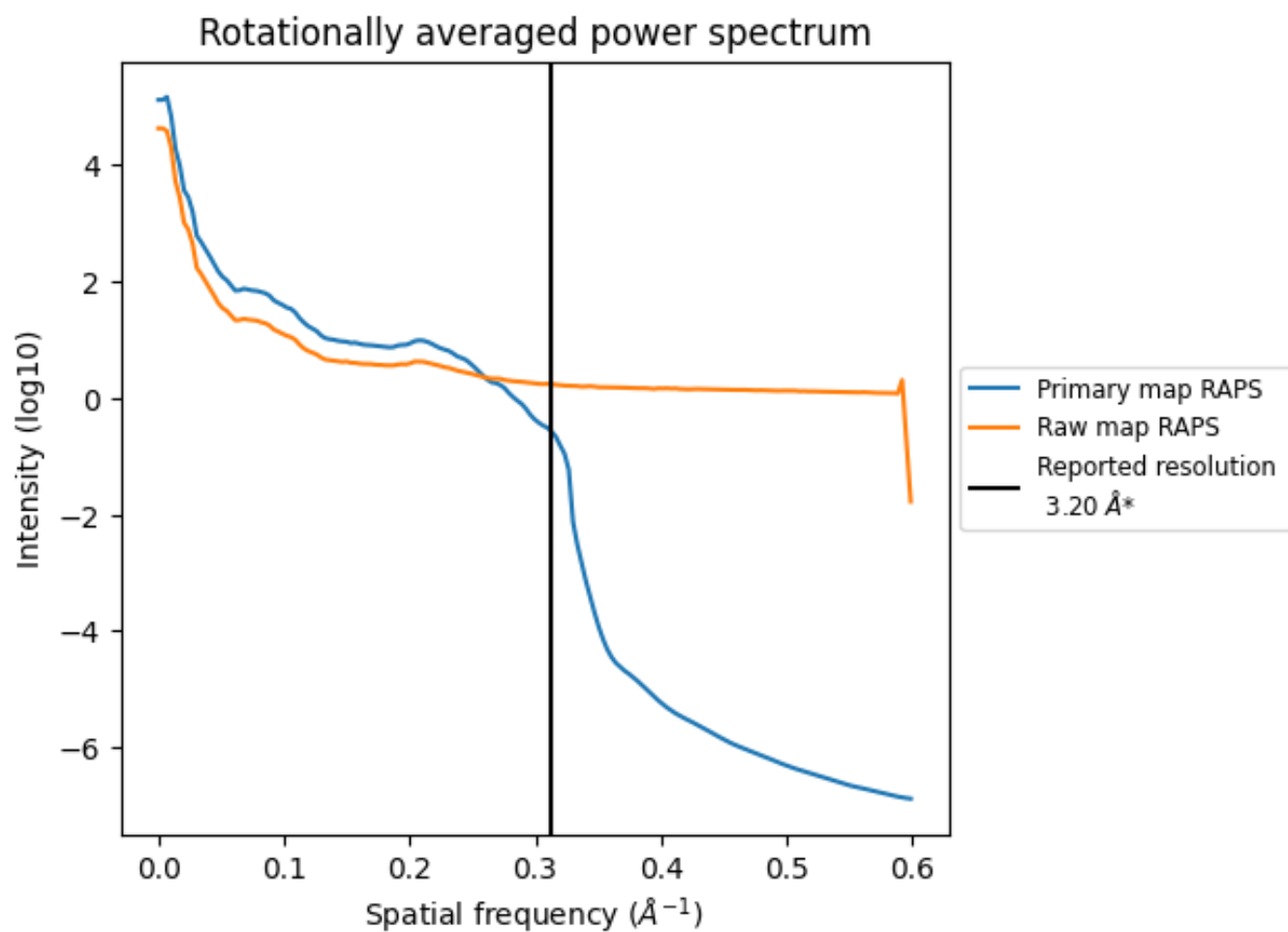
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 166 nm^3 ; this corresponds to an approximate mass of 150 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 [i](#)

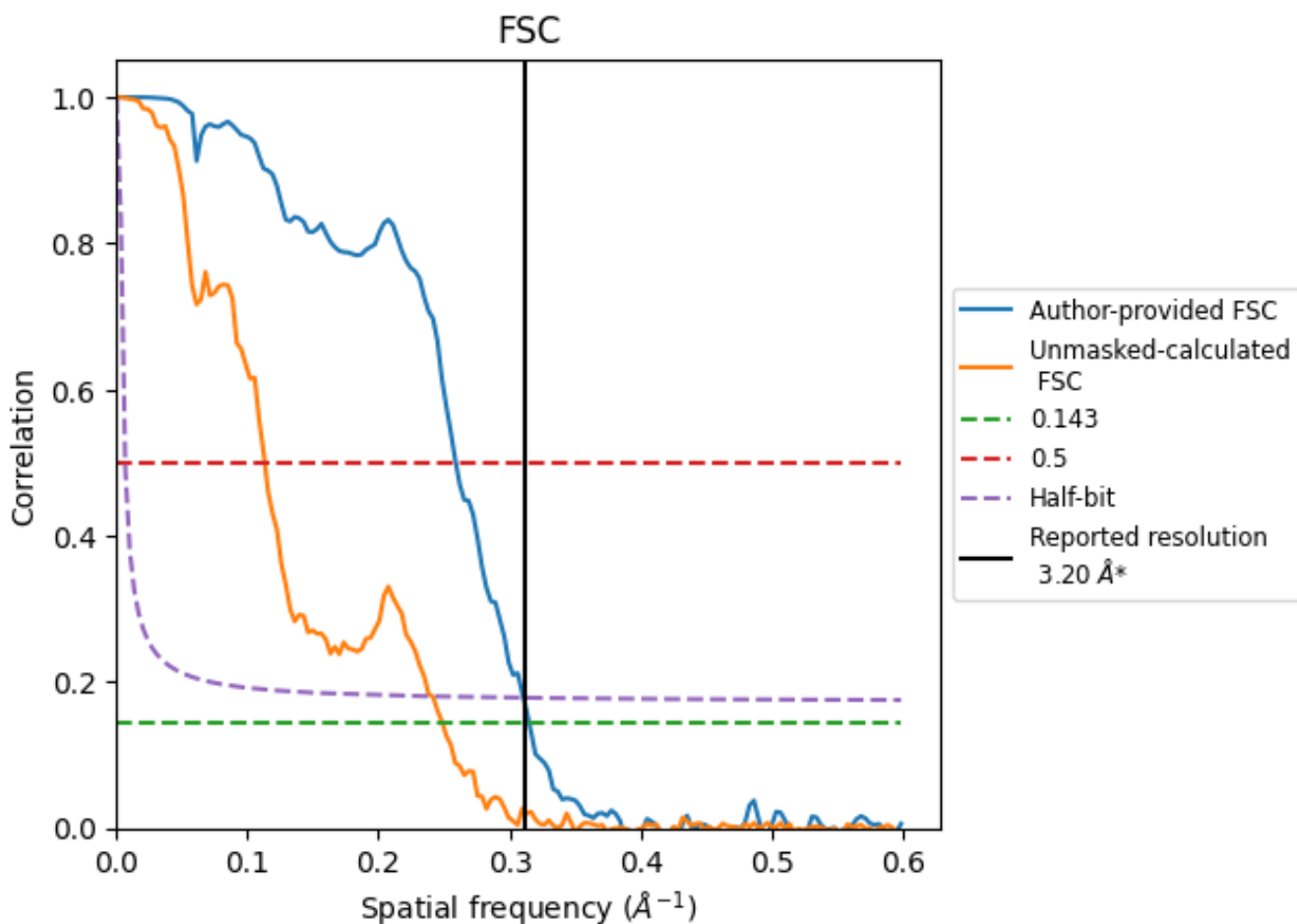


*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

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 0.312\AA^{-1}

8.2 Resolution estimates [i](#)

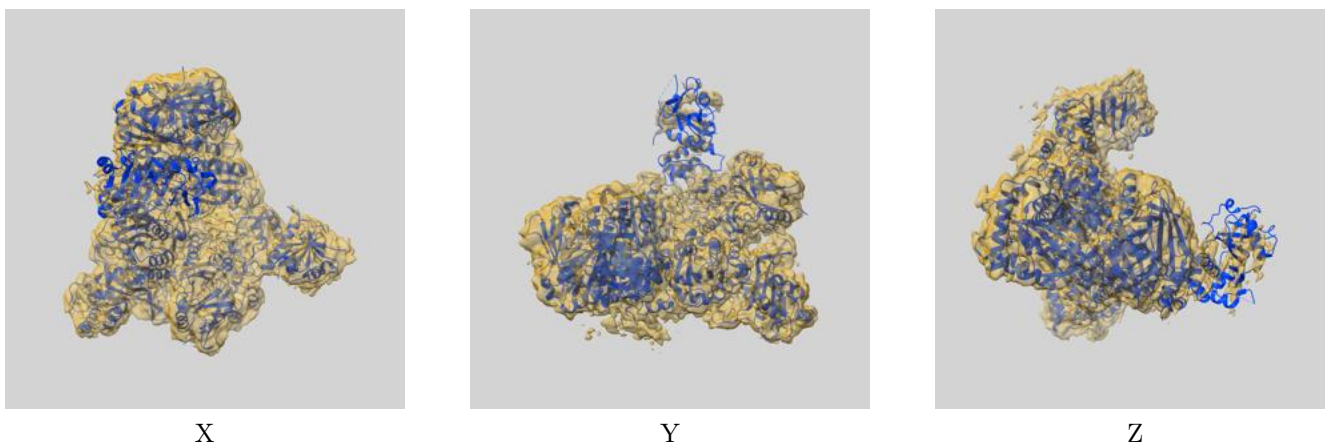
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.17	3.86	3.22
Unmasked-calculated*	4.02	8.83	4.15

*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 4.02 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

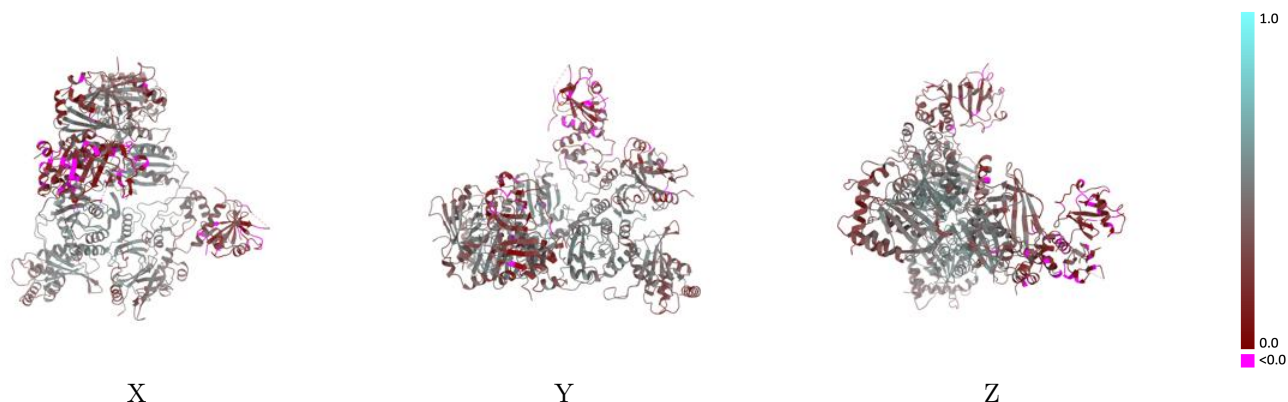
This section contains information regarding the fit between EMDB map EMD-27034 and PDB model 8CX2. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)



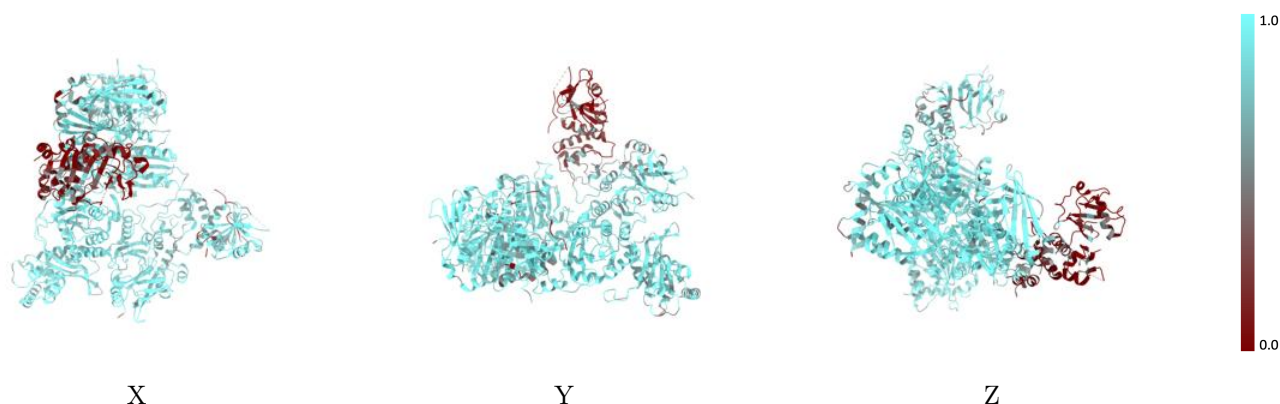
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



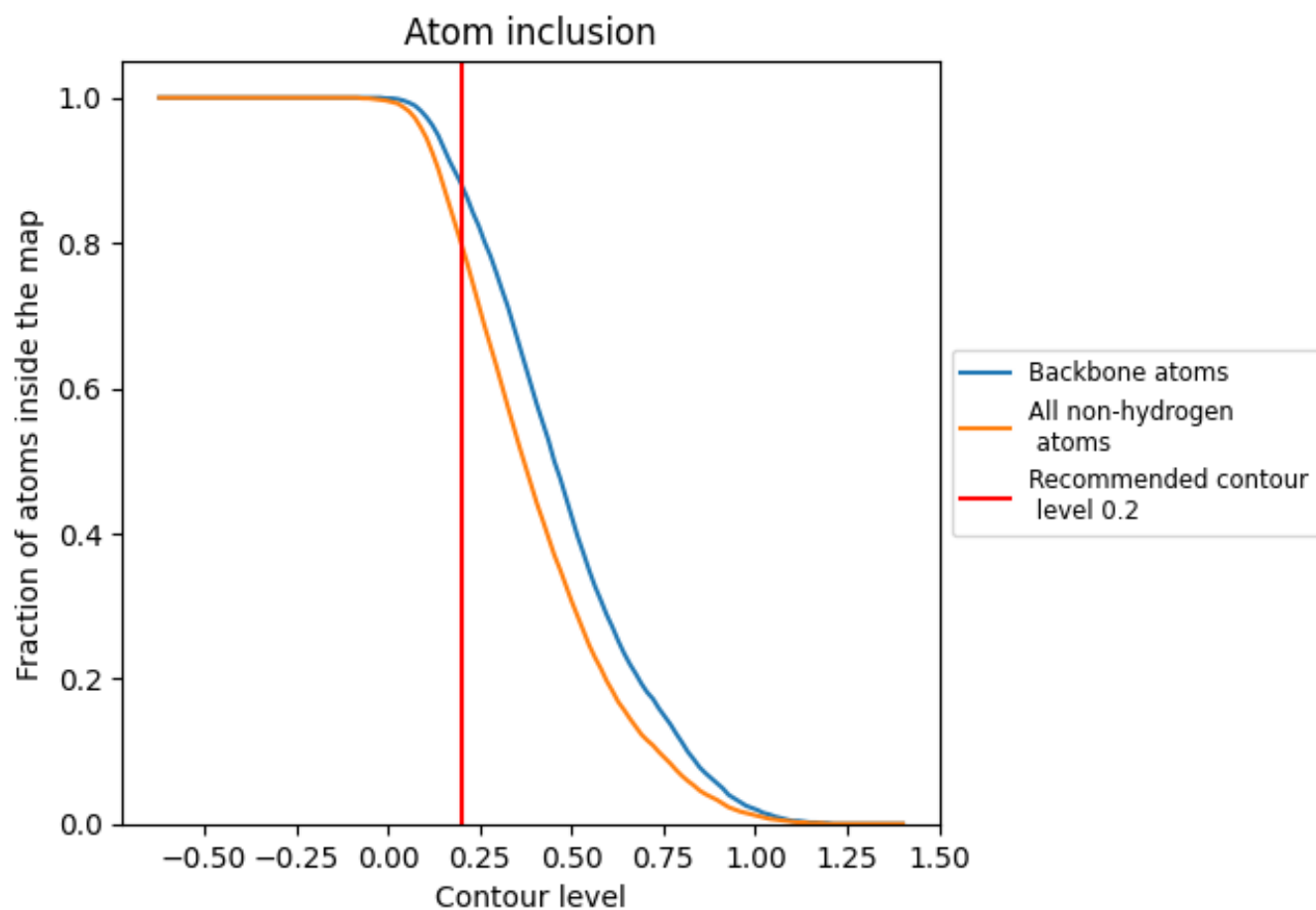
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).



























9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 80% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7990	 0.3750
A	 0.9250	 0.4680
B	 0.9250	 0.4600
C	 0.9190	 0.3940
D	 0.7200	 0.1730
E	 0.7150	 0.2640
F	 0.8890	 0.4270
G	 0.7690	 0.3830
H	 0.7920	 0.2880
I	 0.1270	 0.1020
J	 0.1460	 0.1540
K	 0.9640	 0.5180
L	 0.9680	 0.4730

