



Full wwPDB EM Validation Report ⓘ

Oct 7, 2024 – 05:56 PM EDT

PDB ID : 8V1Q
EMDB ID : EMD-42887
Title : Herpes simplex virus 1 polymerase holoenzyme bound to DNA in both open/closed conformations
Authors : Pan, J.; Abraham, J.; Coen, D.M.; Shankar, S.; Yang, P.; Hogle, J.
Deposited on : 2023-11-21
Resolution : 2.70 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.dev113
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

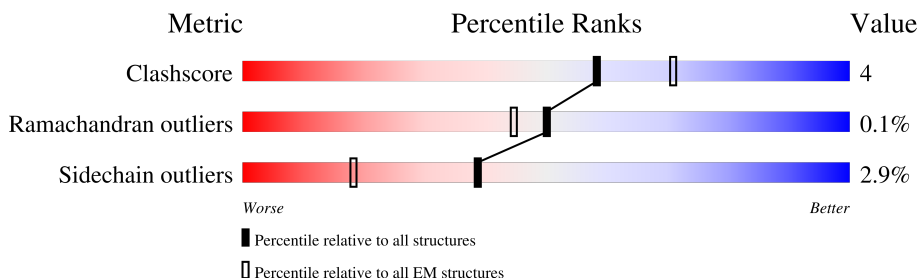
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1199	<div> <div>5%</div> <div>79%</div> <div>12%</div> <div>9%</div> </div>
2	B	349	<div> <div>17%</div> <div>60%</div> <div>16%</div> <div>23%</div> </div>
3	P	32	<div> <div>16%</div> <div>56%</div> <div>16%</div> <div>9%</div> <div>19%</div> </div>
4	T	50	<div> <div>6%</div> <div>32%</div> <div>24%</div> <div>40%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	DOC	P	-1	X	-	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11914 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 polymerase.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1097	8687	5556	1536	1554	41	5	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	HIS	-	expression tag	UNP H9E937
A	38	HIS	-	expression tag	UNP H9E937
A	39	HIS	-	expression tag	UNP H9E937
A	40	HIS	-	expression tag	UNP H9E937
A	41	HIS	-	expression tag	UNP H9E937
A	42	HIS	-	expression tag	UNP H9E937

- Molecule 2 is a protein called DNA polymerase processivity factor.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	270	2077	1321	365	383	8	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	GLY	-	expression tag	UNP A0A181ZFK4
B	-7	PRO	-	expression tag	UNP A0A181ZFK4
B	-6	ILE	-	expression tag	UNP A0A181ZFK4
B	-5	SER	-	expression tag	UNP A0A181ZFK4
B	-4	GLU	-	expression tag	UNP A0A181ZFK4
B	-3	PHE	-	expression tag	UNP A0A181ZFK4
B	-2	GLY	-	expression tag	UNP A0A181ZFK4
B	-1	SER	-	expression tag	UNP A0A181ZFK4
B	0	SER	-	expression tag	UNP A0A181ZFK4
B	1	ARG	-	expression tag	UNP A0A181ZFK4

- Molecule 3 is a DNA chain called Primer DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	P	26	Total	C	N	O	P	0	0
			535	253	101	155	26		

- Molecule 4 is a DNA chain called template DNA.

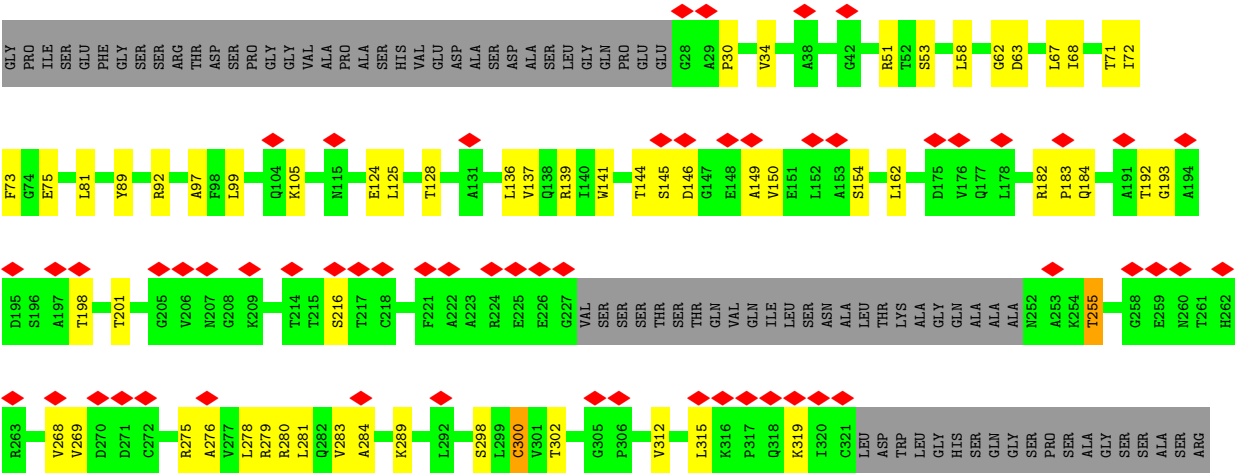
Mol	Chain	Residues	Atoms					AltConf	Trace
4	T	30	Total	C	N	O	P	0	0
			610	289	113	178	30		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

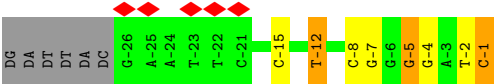
Mol	Chain	Residues	Atoms		AltConf
5	A	2	Total	Mg	0
			2	2	

- Molecule 6 is water.

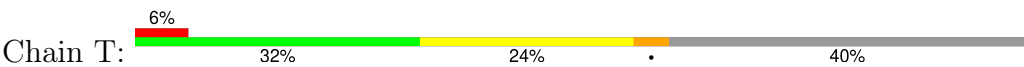
Mol	Chain	Residues	Atoms		AltConf
6	A	3	Total	O	0
			3	3	



• Molecule 3: Primer DNA



• Molecule 4: template DNA



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	343777	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 ($e^-/\text{\AA}^2$)	53.11	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	60606	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.230	Depositor
Minimum map value	-0.123	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.025	Depositor
Map size (\AA)	264.0, 264.0, 264.0	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.825, 0.825, 0.825	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: DOC, MG

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	0.88	4/8890 (0.0%)	0.82	4/12048 (0.0%)
2	B	0.52	0/2115	0.81	0/2877
3	P	1.92	10/580 (1.7%)	1.27	3/894 (0.3%)
4	T	1.89	9/683 (1.3%)	1.49	10/1050 (1.0%)
All	All	0.99	23/12268 (0.2%)	0.90	17/16869 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	B	0	3
3	P	1	0
All	All	1	6

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	-4	DG	C3'-O3'	-8.21	1.33	1.44
1	A	396	CYS	CB-SG	-6.61	1.71	1.82
3	P	-2	DT	C2-N3	-6.61	1.32	1.37
4	T	2	DA	C8-N7	-6.09	1.27	1.31
4	T	7	DC	C3'-O3'	-5.87	1.36	1.44
3	P	-7	DG	N7-C5	-5.69	1.35	1.39
3	P	-7	DG	C8-N7	-5.60	1.27	1.30
1	A	502	TRP	CB-CG	-5.44	1.40	1.50
3	P	-4	DG	C5-C4	-5.40	1.34	1.38
4	T	14	DC	C1'-N1	5.27	1.56	1.49
3	P	-8	DC	N1-C6	-5.25	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	T	2	DA	N7-C5	-5.23	1.36	1.39
4	T	2	DA	C3'-O3'	-5.22	1.37	1.44
3	P	-15	DC	C1'-N1	5.19	1.55	1.49
4	T	3	DT	C3'-O3'	-5.16	1.37	1.44
4	T	2	DA	C6-N1	-5.15	1.31	1.35
1	A	150	TYR	CD2-CE2	-5.13	1.31	1.39
3	P	-2	DT	C5-C6	-5.07	1.30	1.34
3	P	-5	DG	C3'-O3'	-5.05	1.37	1.44
1	A	465	TYR	CD2-CE2	-5.03	1.31	1.39
3	P	-7	DG	C6-N1	-5.02	1.36	1.39
4	T	9	DG	C5-C4	-5.02	1.34	1.38
4	T	1	DG	C3'-O3'	-5.02	1.37	1.44

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	T	-1	DC	O3'-P-O5'	-15.08	75.34	104.00
4	T	-1	DC	P-O3'-C3'	12.04	134.15	119.70
4	T	-1	DC	OP1-P-O3'	-11.66	79.54	105.20
4	T	-3	DC	P-O3'-C3'	10.15	131.88	119.70
4	T	2	DA	P-O3'-C3'	9.55	131.16	119.70
4	T	17	DG	P-O3'-C3'	9.12	130.64	119.70
1	A	867	PRO	CA-N-CD	-8.24	99.96	111.50
3	P	-4	DG	O4'-C1'-N9	7.75	113.43	108.00
4	T	0	DA	P-O3'-C3'	6.77	127.82	119.70
4	T	0	DA	OP1-P-OP2	6.52	129.38	119.60
1	A	1182	LYS	CG-CD-CE	5.75	129.16	111.90
1	A	1182	LYS	CD-CE-NZ	5.72	124.86	111.70
4	T	3	DT	P-O3'-C3'	5.67	126.51	119.70
1	A	990	LEU	CA-CB-CG	5.59	128.17	115.30
3	P	-4	DG	C1'-O4'-C4'	-5.54	104.56	110.10
3	P	-12	DT	N3-C4-O4	5.28	123.07	119.90
4	T	-1	DC	OP2-P-O3'	-5.15	93.88	105.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	P	-1	DOC	C4'

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	176	THR	Peptide
1	A	259	TYR	Peptide
1	A	956	ASP	Peptide
2	B	150	VAL	Peptide
2	B	319	LYS	Mainchain
2	B	53	SER	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8687	0	8687	74	0
2	B	2077	0	2125	29	0
3	P	535	0	291	3	0
4	T	610	0	336	8	0
5	A	2	0	0	0	0
6	A	3	0	0	1	0
All	All	11914	0	11439	105	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 (105) 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:511:LYS:CE	4:T:0:DA:OP1	2.20	0.89
1:A:311:VAL:HG22	1:A:352:ILE:HD11	1.56	0.84
1:A:1215:GLY:O	1:A:1223:THR:OG1	1.97	0.81
2:B:192:THR:O	2:B:275:ARG:NH2	2.15	0.80
1:A:511:LYS:HE3	4:T:0:DA:OP1	1.83	0.78
1:A:346:THR:OG1	1:A:348:ASP:OD1	2.05	0.72
1:A:977:LEU:O	1:A:984:SER:OG	2.08	0.70
2:B:280:ARG:NE	4:T:19:DT:OP1	2.24	0.69
1:A:1229:ARG:O	1:A:1233:THR:OG1	2.05	0.69
2:B:289:LYS:NZ	2:B:302:THR:OG1	2.25	0.69
1:A:1194:GLU:HG3	2:B:162:LEU:HD23	1.75	0.68
1:A:616:GLY:O	6:A:1501:HOH:O	2.11	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:534:LYS:NZ	1:A:958:VAL:O	2.18	0.67
1:A:244:SER:OG	1:A:245:ALA:N	2.27	0.67
1:A:216:LEU:O	1:A:227[A]:ARG:NH1	2.29	0.66
2:B:72:ILE:HD12	2:B:281:LEU:HD21	1.78	0.66
1:A:239:SER:O	1:A:277:ARG:NH2	2.27	0.65
1:A:325:THR:O	1:A:349:ASN:ND2	2.33	0.60
1:A:187:LEU:HD23	1:A:360:PRO:O	2.04	0.58
1:A:1054:VAL:HG21	1:A:1074:TYR:HB3	1.85	0.57
1:A:212:VAL:HG12	1:A:290:ILE:HD13	1.85	0.57
2:B:34:VAL:HG22	2:B:124:GLU:HG2	1.87	0.57
1:A:1221:GLU:OE1	1:A:1224:ARG:NH2	2.38	0.57
1:A:873:ARG:NH2	1:A:876:GLY:O	2.38	0.56
1:A:884:TYR:OH	1:A:938:LYS:O	2.17	0.56
1:A:732:CYS:SG	1:A:733:PHE:N	2.78	0.56
1:A:1034:THR:HG23	1:A:1071:ARG:HH21	1.71	0.56
2:B:125:LEU:HG	2:B:136:LEU:HD11	1.88	0.55
2:B:30:PRO:HD2	2:B:128:THR:HG22	1.88	0.55
2:B:30:PRO:CD	2:B:128:THR:HG22	2.37	0.55
1:A:598:LEU:HD13	1:A:613:ILE:HG12	1.89	0.55
2:B:72:ILE:CD1	2:B:281:LEU:HD21	2.36	0.54
1:A:536:SER:HB3	1:A:1034:THR:HG21	1.88	0.54
1:A:193:ARG:NH1	1:A:458:GLY:O	2.42	0.53
1:A:310:PHE:HA	1:A:352:ILE:HD13	1.91	0.53
1:A:1054:VAL:HG21	1:A:1074:TYR:CB	2.39	0.52
2:B:97:ALA:HB3	2:B:162:LEU:HD12	1.90	0.52
2:B:141:TRP:HD1	2:B:149:ALA:HB1	1.72	0.52
1:A:89:GLU:OE1	1:A:89:GLU:N	2.42	0.52
1:A:757:VAL:HG21	1:A:762:LEU:HD12	1.90	0.52
1:A:117:VAL:HG12	1:A:117:VAL:O	2.10	0.52
1:A:511:LYS:NZ	4:T:0:DA:OP1	2.42	0.52
1:A:1146:PRO:O	1:A:1150:ILE:HD12	2.10	0.52
1:A:700:ARG:HH22	1:A:846:LEU:HD11	1.74	0.52
1:A:73:GLU:OE2	1:A:98:HIS:ND1	2.41	0.52
2:B:62:GLY:O	2:B:89:TYR:OH	2.18	0.51
1:A:818:TYR:O	1:A:821:THR:OG1	2.26	0.51
1:A:389:ASP:O	1:A:478:LYS:NZ	2.27	0.51
2:B:193:GLY:HA3	2:B:198:THR:HG21	1.93	0.51
2:B:269:VAL:HG22	2:B:315:LEU:HD22	1.91	0.50
1:A:363:LYS:NZ	1:A:400:ASP:OD2	2.40	0.50
2:B:182:ARG:HE	2:B:284:ALA:HB1	1.76	0.49
1:A:281:TYR:CE2	1:A:285:ASN:ND2	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:68:ILE:HD11	2:B:81:LEU:HD11	1.94	0.49
1:A:273:VAL:HG21	1:A:279:LEU:HB2	1.95	0.49
1:A:1048:ARG:NH1	1:A:1141:GLU:OE2	2.46	0.48
1:A:819:GLY:O	1:A:823:VAL:HG23	2.13	0.48
2:B:105:LYS:NZ	3:P:-12:DT:OP1	2.38	0.48
1:A:640:GLN:NE2	4:T:-1:DC:N3	2.49	0.48
2:B:58:LEU:HD11	2:B:162:LEU:HD11	1.96	0.48
1:A:251:GLU:OE2	1:A:270:ARG:NH2	2.41	0.48
2:B:97:ALA:CB	2:B:162:LEU:HD12	2.42	0.48
1:A:552:LYS:NZ	1:A:581:ASP:OD1	2.28	0.48
1:A:148:HIS:O	1:A:186:GLY:HA3	2.14	0.47
1:A:232:LEU:HD11	1:A:278:VAL:HG13	1.95	0.47
2:B:201:THR:HG22	2:B:268:VAL:HG23	1.97	0.47
3:P:-1:DOC:H6	3:P:-1:DOC:H4'	1.97	0.47
4:T:24:DT:H2''	4:T:25:DT:H71	1.96	0.47
1:A:194:VAL:HG23	1:A:341:VAL:HG23	1.96	0.47
1:A:187:LEU:HD22	1:A:362:TYR:CD1	2.50	0.46
1:A:640:GLN:HA	1:A:643:PHE:HB3	1.96	0.46
2:B:278:LEU:CD2	2:B:283:VAL:HG21	2.46	0.46
1:A:298:ASP:OD1	1:A:298:ASP:N	2.46	0.45
1:A:557:TYR:HA	1:A:560:ILE:HD12	1.99	0.45
1:A:628:LEU:HD21	1:A:776:ILE:HB	1.98	0.45
1:A:698:GLY:HA2	1:A:838:THR:HG21	1.97	0.45
1:A:257:ASP:OD1	1:A:258:VAL:N	2.49	0.44
1:A:736:LEU:HD13	1:A:764:PHE:CZ	2.52	0.44
1:A:364:LEU:HB2	1:A:401:LEU:HD13	2.00	0.44
1:A:699:ALA:HB2	1:A:887:THR:HG23	2.00	0.44
1:A:1206:LEU:HA	1:A:1209:ALA:HB3	2.00	0.43
1:A:1214:VAL:O	1:A:1223:THR:HG23	2.18	0.43
1:A:76:PHE:CD2	1:A:99:LEU:HD22	2.53	0.43
1:A:180:THR:HG21	1:A:299:ALA:CB	2.49	0.43
2:B:63:ASP:N	2:B:63:ASP:OD1	2.50	0.43
4:T:23:DA:H2'	4:T:24:DT:H72	2.01	0.43
1:A:785:ARG:HB2	1:A:810:ILE:HG21	2.00	0.43
1:A:476:LEU:HD22	1:A:487:LEU:HD13	1.99	0.43
1:A:511:LYS:HE2	4:T:0:DA:OP1	2.11	0.43
1:A:206:TYR:HB2	1:A:293:TYR:HB2	2.01	0.42
1:A:233:ARG:NH1	1:A:243:ILE:O	2.51	0.42
1:A:849:ARG:HA	1:A:882:ILE:HD13	2.01	0.42
2:B:300:CYS:HB2	2:B:312:VAL:HG22	2.01	0.42
1:A:723:PRO:HB3	1:A:778:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:PRO:HG2	1:A:341:VAL:HG11	2.01	0.42
2:B:99:LEU:HG	2:B:162:LEU:HD21	2.01	0.42
1:A:999:LEU:HD21	1:A:1183:ILE:HG23	2.02	0.41
1:A:872:MET:CE	1:A:902:LEU:HD13	2.50	0.41
1:A:1041:PRO:HB3	1:A:1052:LEU:HD22	2.03	0.41
2:B:68:ILE:CD1	2:B:81:LEU:HD11	2.50	0.41
2:B:73:PHE:CE2	2:B:276:ALA:HB1	2.56	0.41
2:B:30:PRO:HD3	2:B:128:THR:HG22	2.03	0.41
1:A:464:GLY:HA2	1:A:585:VAL:HG11	2.03	0.41
1:A:692:ARG:NH2	3:P:-5:DG:N7	2.68	0.41
2:B:184:GLN:NE2	2:B:255:THR:O	2.54	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	1096/1199 (91%)	1091 (100%)	5 (0%)	0	100	100
2	B	266/349 (76%)	265 (100%)	0	1 (0%)	30	55
All	All	1362/1548 (88%)	1356 (100%)	5 (0%)	1 (0%)	50	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	ARG

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	905/977 (93%)	888 (98%)	17 (2%)	52	79
2	B	230/289 (80%)	214 (93%)	16 (7%)	12	31
All	All	1135/1266 (90%)	1102 (97%)	33 (3%)	39	67

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	239	SER
1	A	264	ARG
1	A	298	ASP
1	A	344	ASN
1	A	390	LEU
1	A	539	LYS
1	A	618	GLN
1	A	739	ARG
1	A	741	ASP
1	A	747	GLU
1	A	872	MET
1	A	907	ASP
1	A	1008	GLN
1	A	1019	ARG
1	A	1062	ARG
1	A	1170	CYS
1	A	1189	LYS
2	B	67	LEU
2	B	71	THR
2	B	75	GLU
2	B	92	ARG
2	B	137	VAL
2	B	139	ARG
2	B	144	THR
2	B	145	SER
2	B	146	ASP
2	B	154	SER
2	B	183	PRO
2	B	216	SER
2	B	255	THR
2	B	279	ARG
2	B	298	SER
2	B	300	CYS

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DOC	P	-1	3,4	16,19,20	2.32	8 (50%)	20,26,29	2.62	4 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DOC	P	-1	3,4	1/1/3/3	2/7/18/19	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	-1	DOC	C5-C4	-3.86	1.34	1.42
3	P	-1	DOC	O2-C2	-3.61	1.17	1.23
3	P	-1	DOC	C4-N3	-3.57	1.27	1.34
3	P	-1	DOC	C3'-C2'	-3.16	1.45	1.54
3	P	-1	DOC	C2-N1	-2.97	1.33	1.40
3	P	-1	DOC	C4-N4	2.88	1.40	1.33
3	P	-1	DOC	C6-N1	-2.51	1.32	1.38
3	P	-1	DOC	O5'-C5'	-2.13	1.37	1.44

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	-1	DOC	C4'-O4'-C1'	-10.02	100.35	109.81
3	P	-1	DOC	O4'-C4'-C5'	2.96	114.67	109.34
3	P	-1	DOC	O4'-C1'-N1	2.89	112.99	107.86
3	P	-1	DOC	O2-C2-N3	-2.71	118.06	122.33

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	P	-1	DOC	C4'

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	-1	DOC	C3'-C4'-C5'-O5'
3	P	-1	DOC	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	-1	DOC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

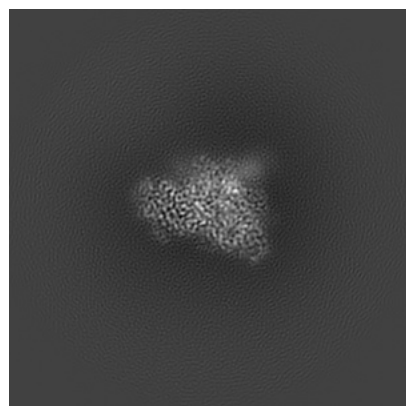
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-42887. These allow visual inspection of the internal detail of the map and identification of artifacts.

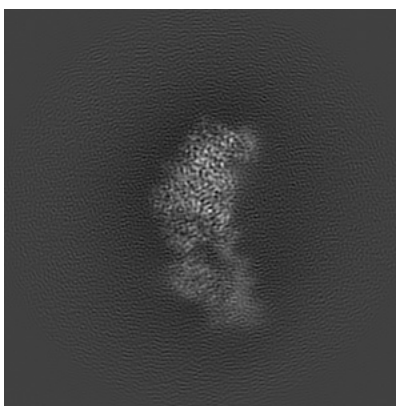
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

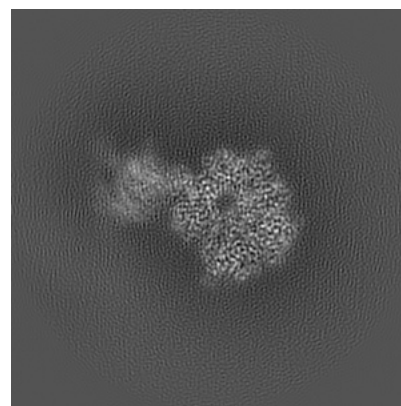
6.1.1 Primary map



X

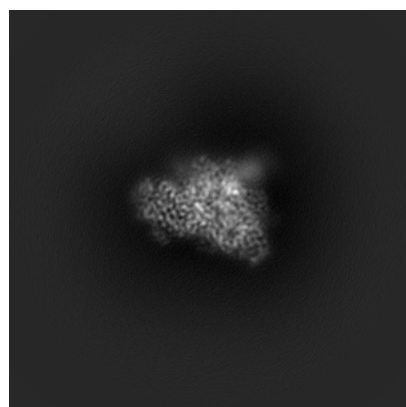


Y

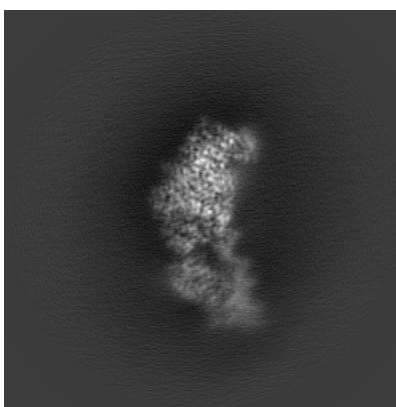


Z

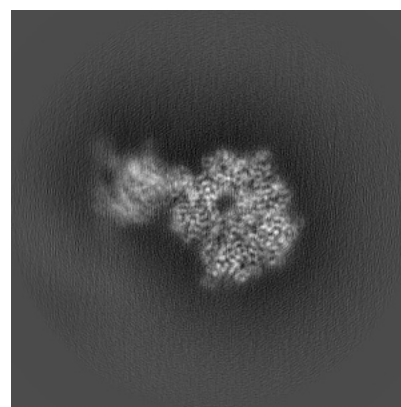
6.1.2 Raw map



X



Y

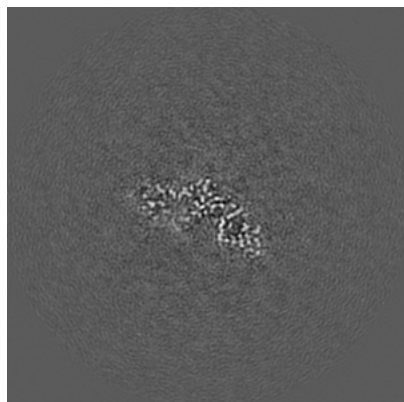


Z

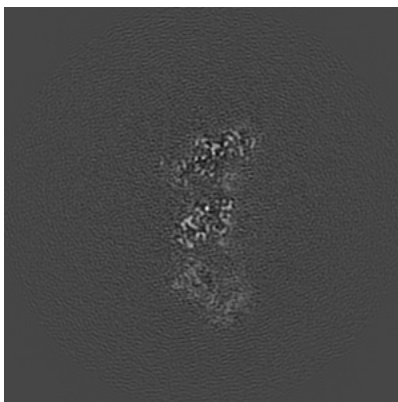
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

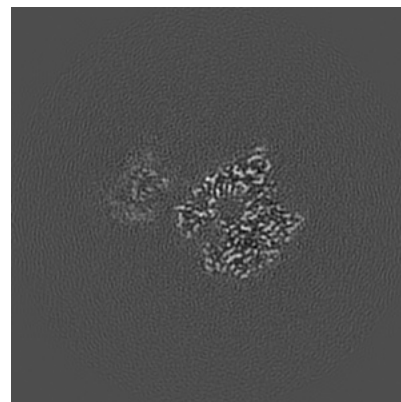
6.2.1 Primary map



X Index: 160

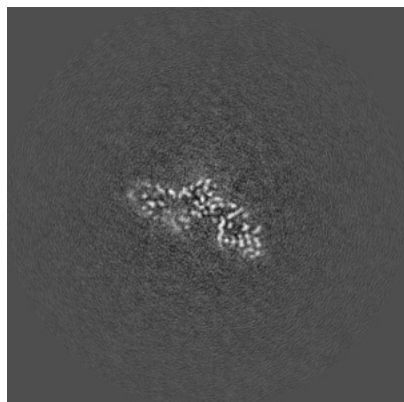


Y Index: 160

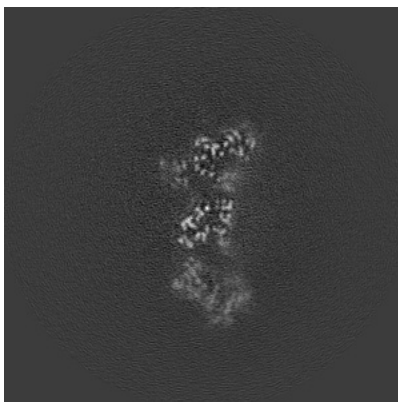


Z Index: 160

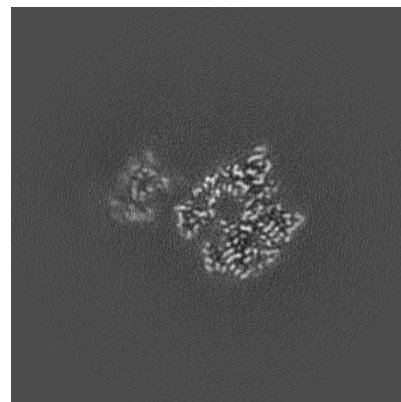
6.2.2 Raw map



X Index: 160



Y Index: 160

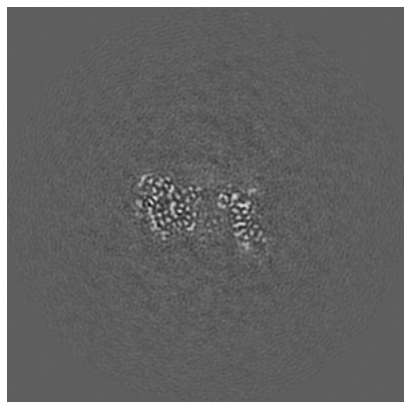


Z Index: 160

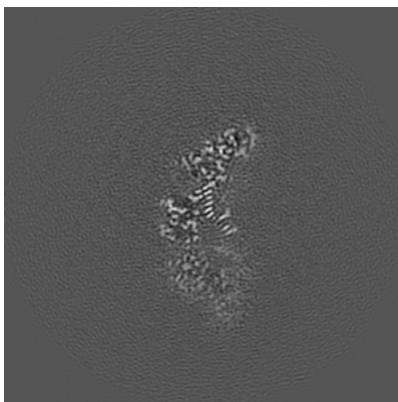
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

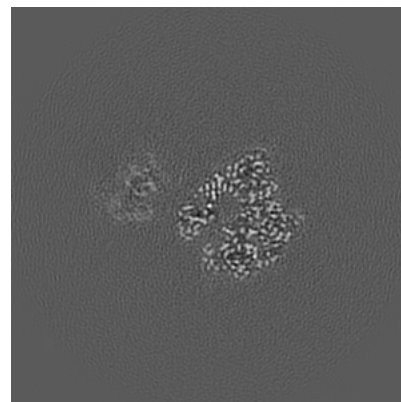
6.3.1 Primary map



X Index: 179

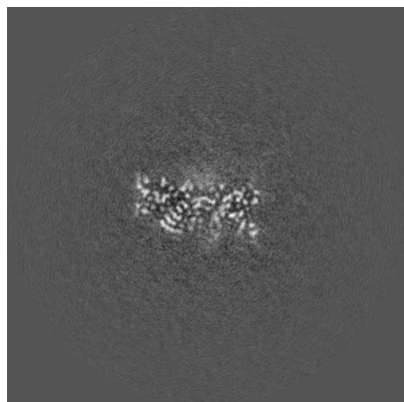


Y Index: 175

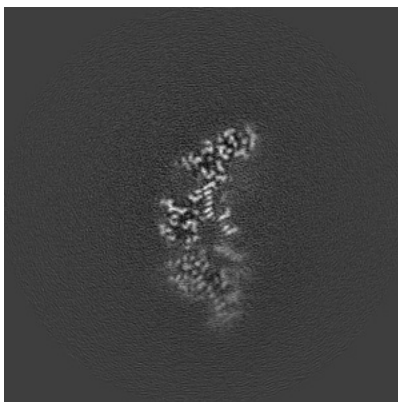


Z Index: 162

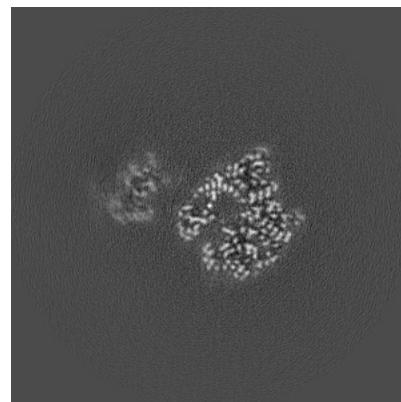
6.3.2 Raw map



X Index: 186



Y Index: 175

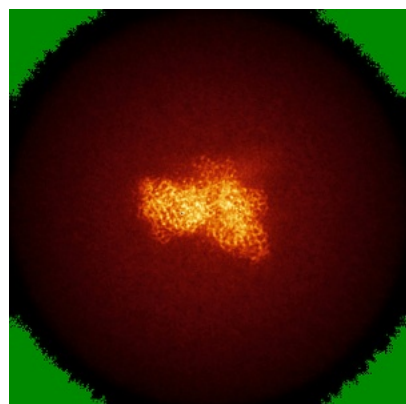


Z Index: 162

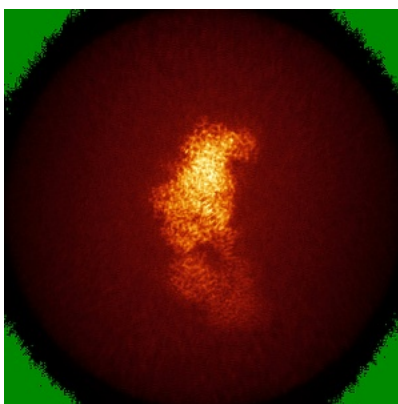
The images above show the largest variance slices of the map in three orthogonal directions.

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

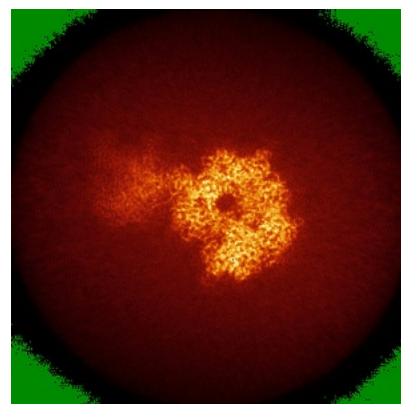
6.4.1 Primary map



X

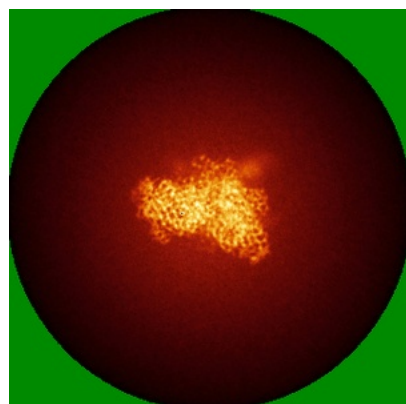


Y

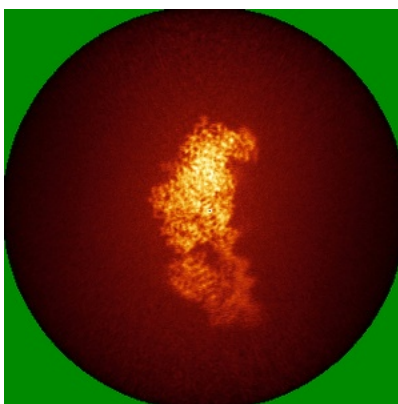


Z

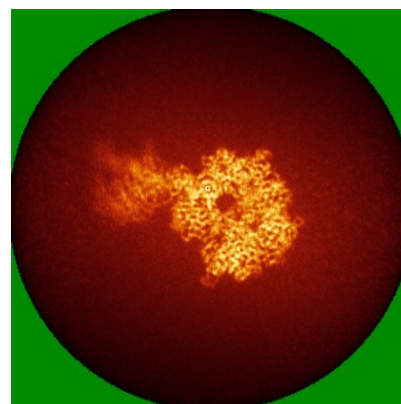
6.4.2 Raw map



X



Y

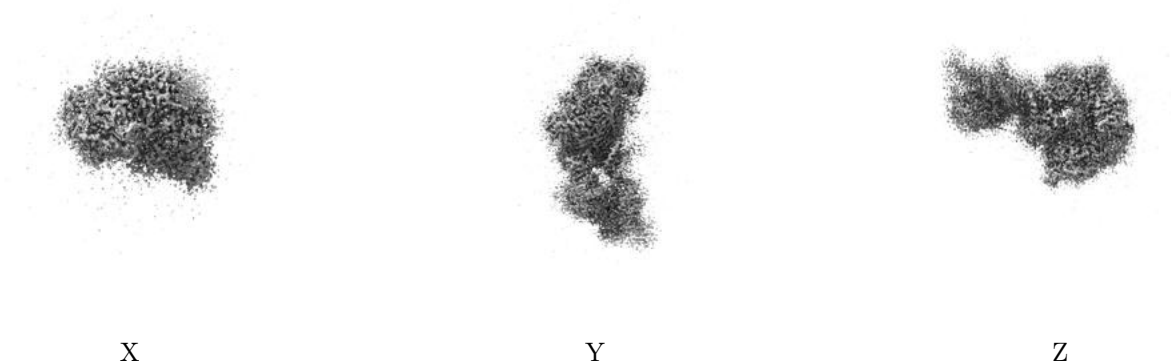


Z

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.

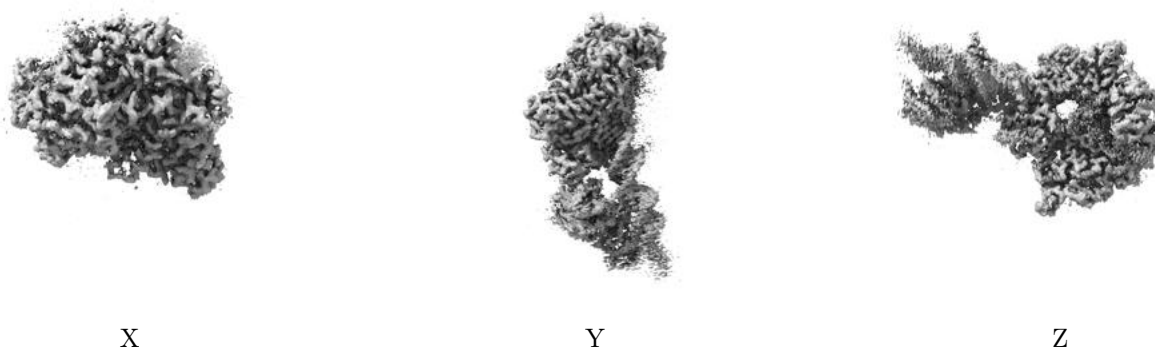
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. 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.

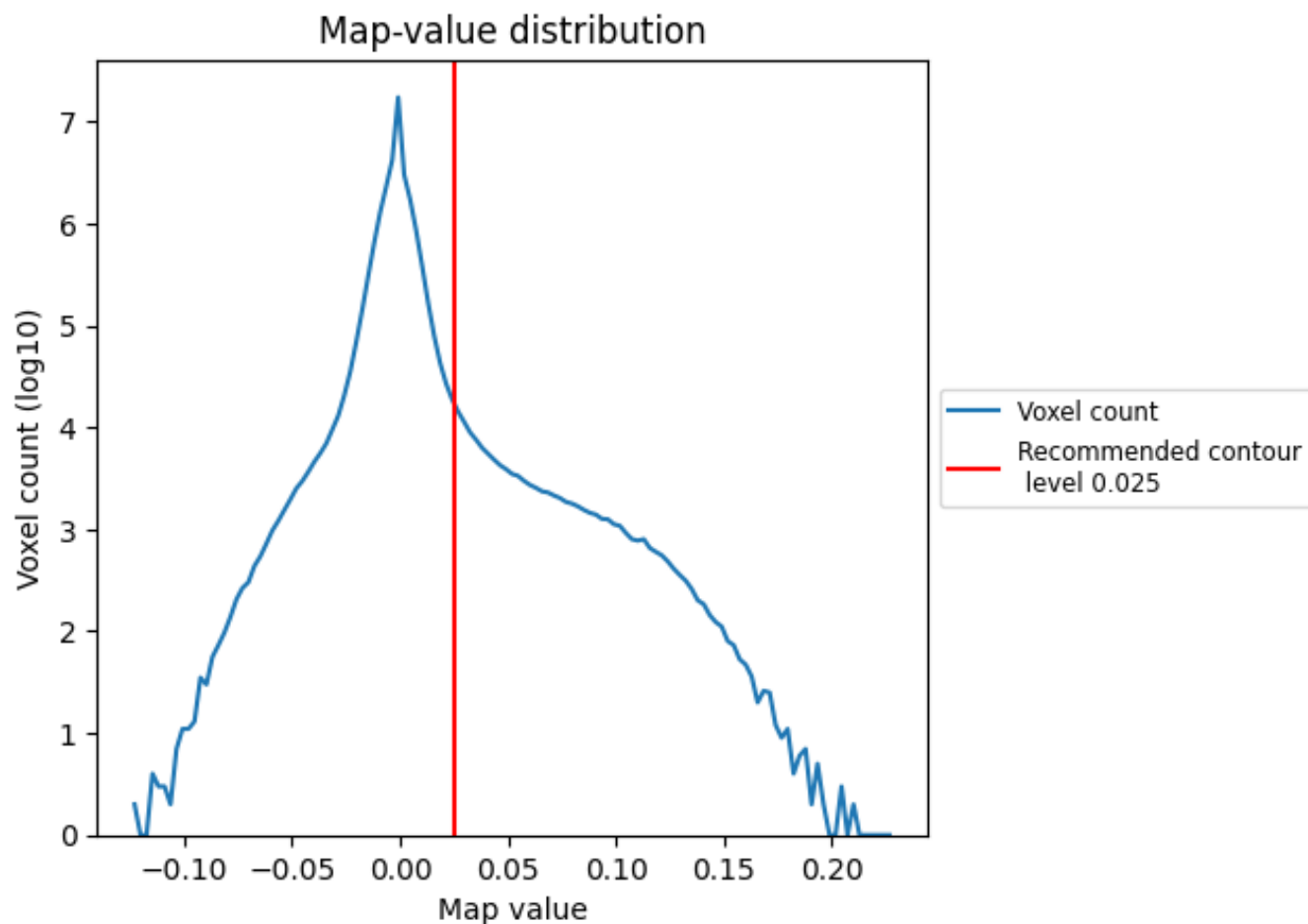
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

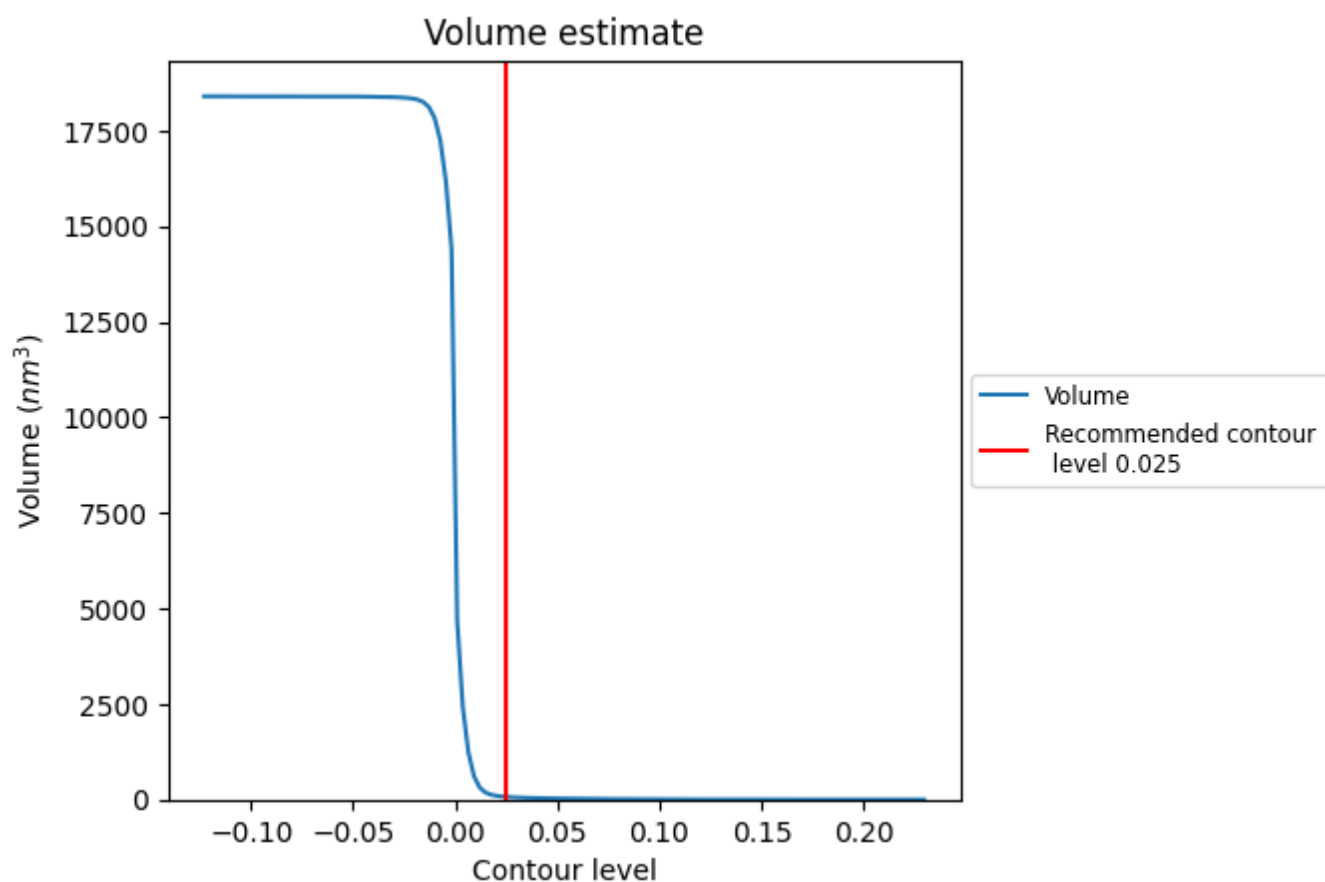
This section contains the results of statistical analysis of the map.

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.

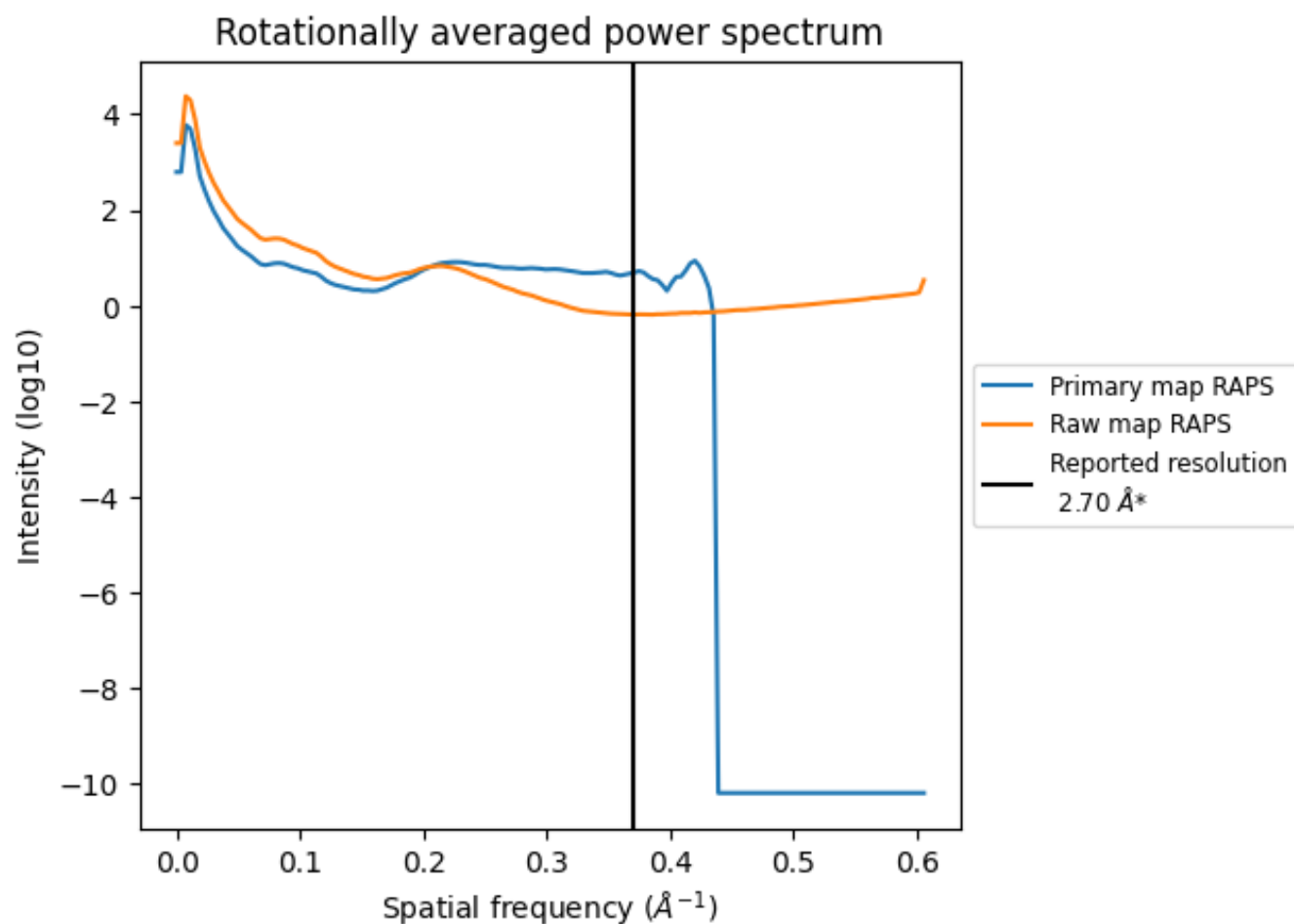
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 71 nm³; this corresponds to an approximate mass of 64 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 ⓘ

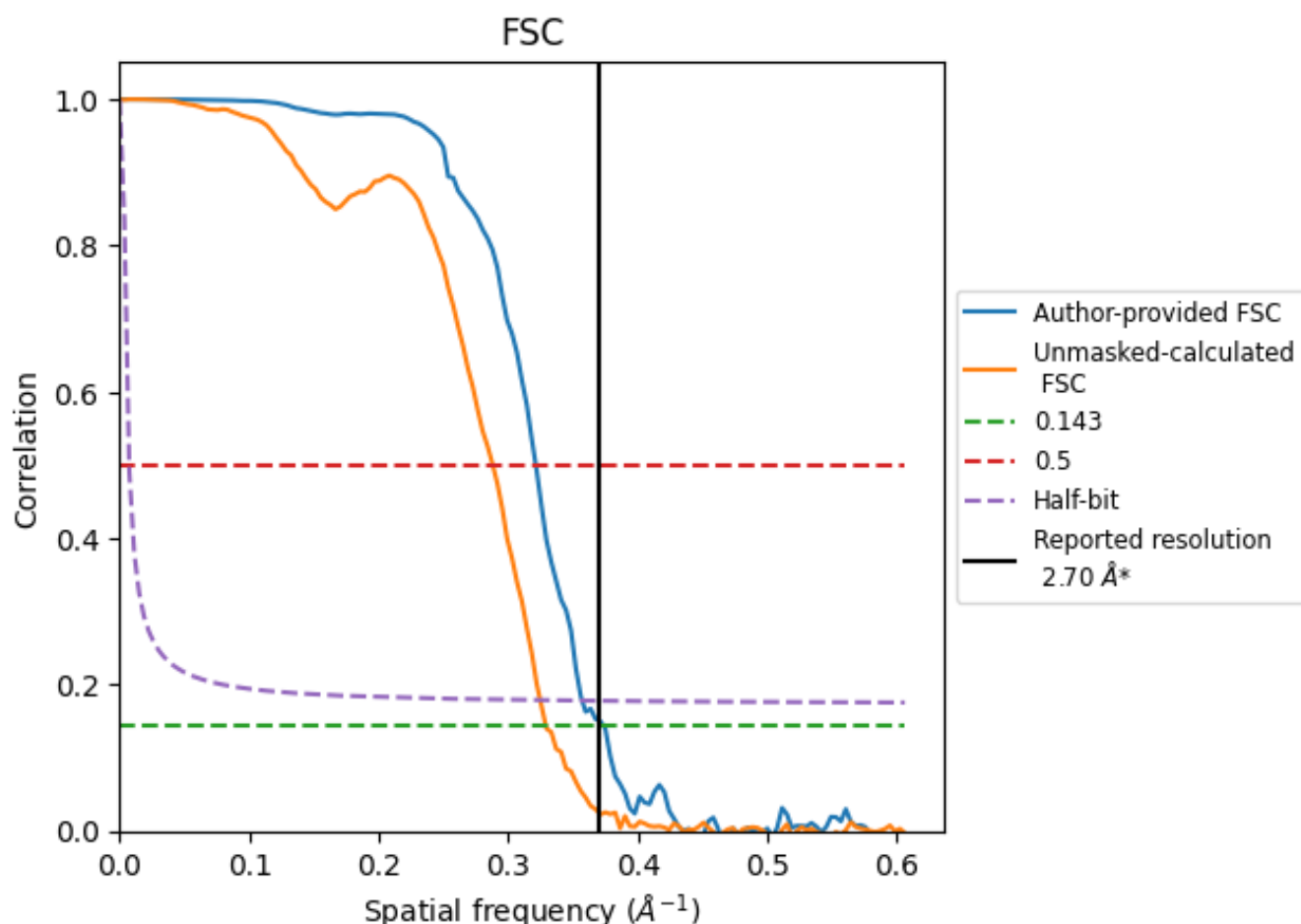


*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

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 0.370 Å⁻¹

8.2 Resolution estimates [i](#)

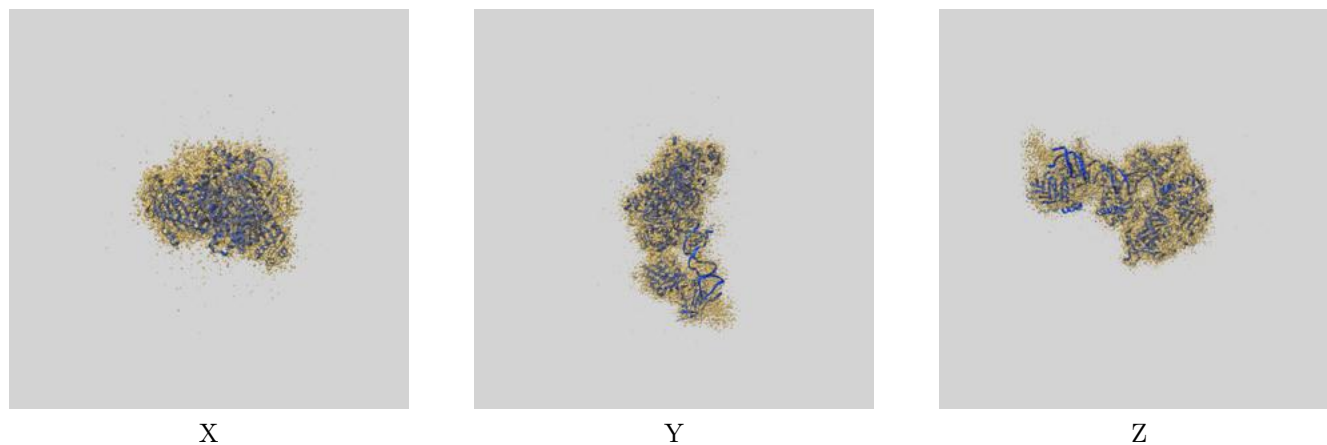
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	2.68	3.11	2.80
Unmasked-calculated*	3.04	3.47	3.08

*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 3.04 differs from the reported value 2.7 by more than 10 %

9 Map-model fit [i](#)

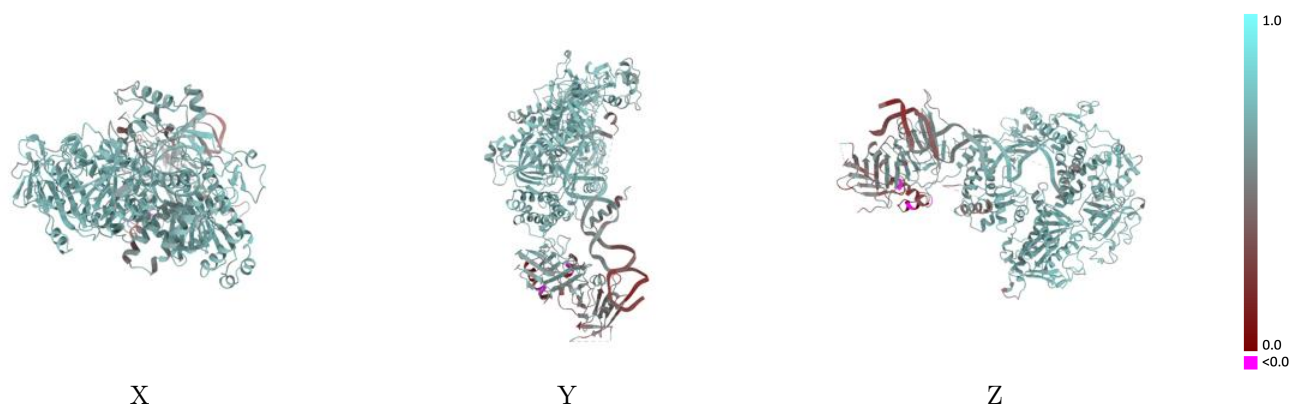
This section contains information regarding the fit between EMDB map EMD-42887 and PDB model 8V1Q. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

9.1 Map-model overlay [i](#)



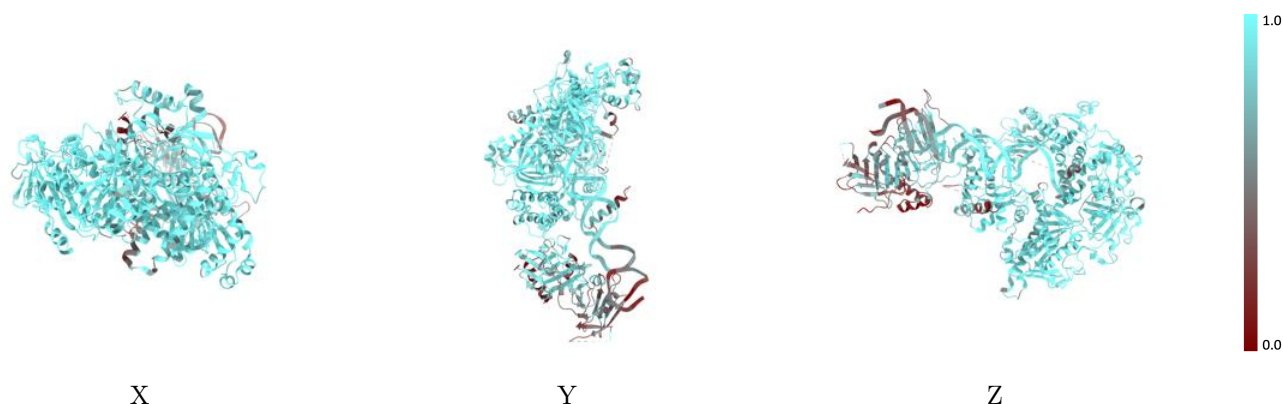
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



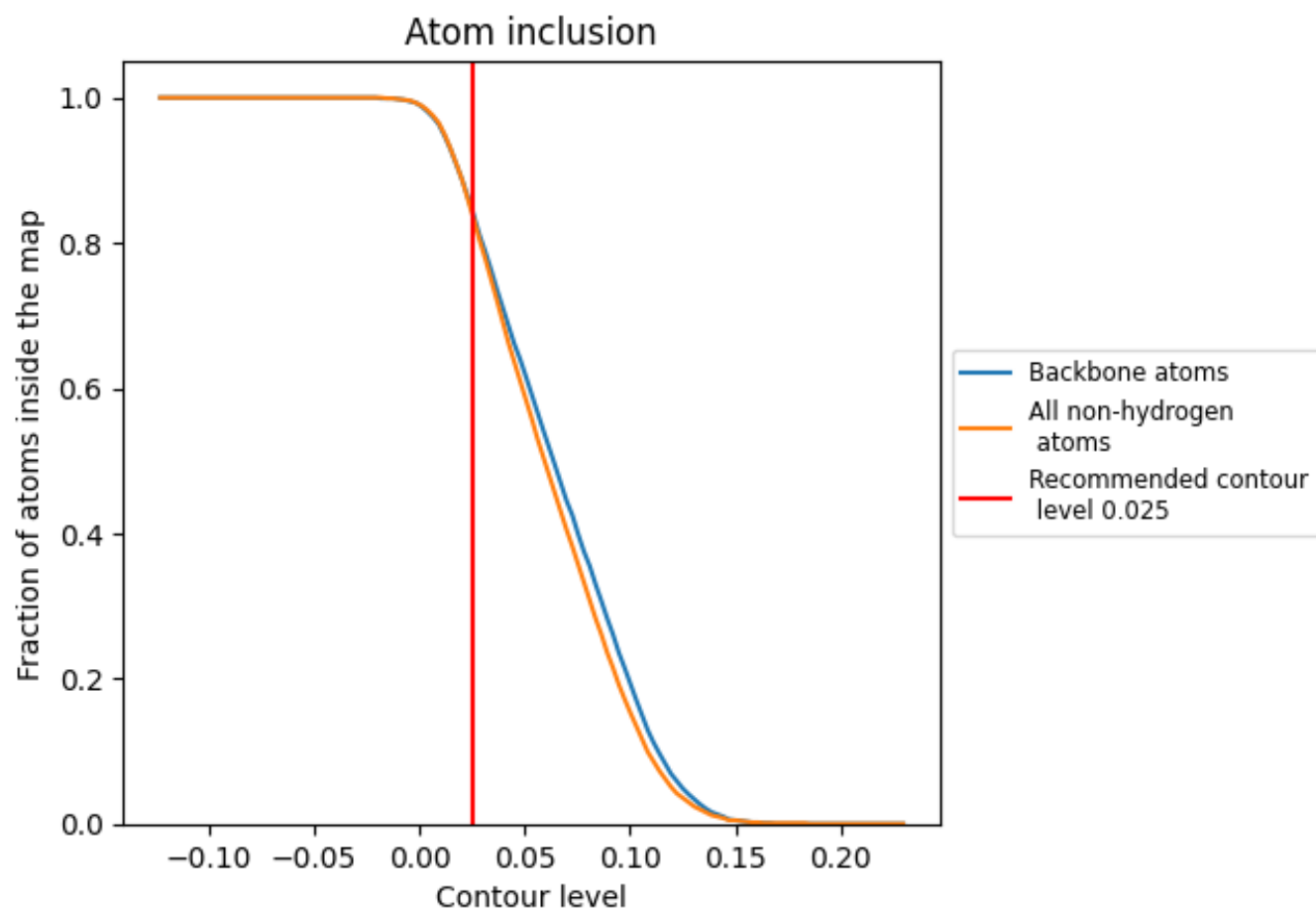
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 84% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8410	<div></div> 0.6030
A	<div></div> 0.9060	<div></div> 0.6390
B	<div></div> 0.6350	<div></div> 0.5070
P	<div></div> 0.6860	<div></div> 0.4730
T	<div></div> 0.7750	<div></div> 0.5240

