



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 01:40 AM UTC

PDB ID : 7OHA / pdb_00007oha
EMDB ID : EMD-12898
Title : nucleosome with TBP and TFIIA bound at SHL +2
Authors : Wang, H.; Cramer, P.
Deposited on : 2021-05-09
Resolution : 2.90 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:

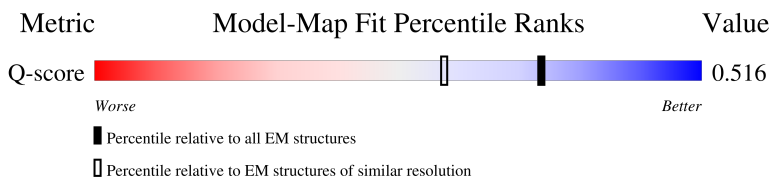
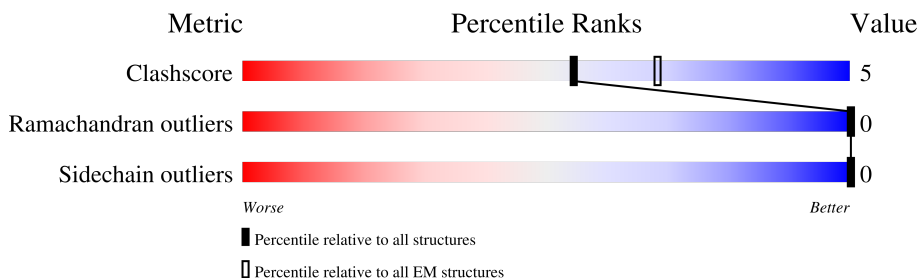
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 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.






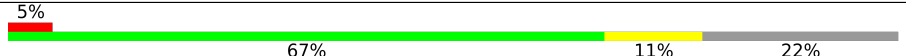
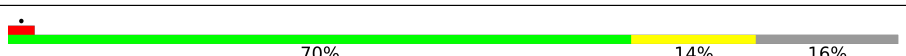
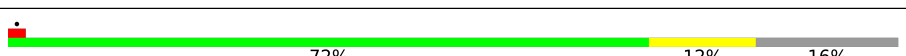
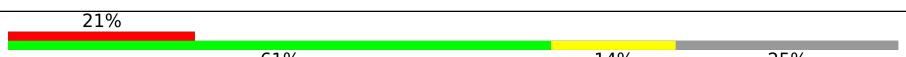
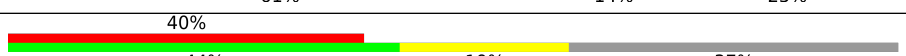

Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13054 (2.40 - 3.40)

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	135	 61% 10% 28%
1	E	135	 64% 8% 27%
2	B	102	 67% 14% 20%
2	F	102	 74% 5% 22%

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Mol	Chain	Length	Quality of chain
3	C	129	 77% 8% 16%
3	G	129	 72% 10% 18%
4	D	122	 70% 10% 20%
4	H	122	 5% 67% 11% 22%
5	I	145	 70% 14% 16%
6	J	145	 72% 12% 16%
7	K	240	 21% 61% 14% 25%
8	L	171	 40% 44% 19% 37%
9	M	122	 47% 64% 23% 12%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 14215 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 Histone H3.2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	97	802	506	155	138	3	0	0
1	E	98	811	512	157	139	3	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	GLY	conflict	UNP P84233
E	102	ALA	GLY	conflict	UNP P84233

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	82	653	412	127	113	1	0	0
2	F	80	638	401	125	111	1	0	0

- Molecule 3 is a protein called Histone H2A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	C	109	843	531	167	145	0	0
3	G	106	818	516	160	142	0	0

- Molecule 4 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	97	767	481	142	142	2	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	H	95	745	469	134	140	2	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	29	THR	SER	conflict	UNP P02281
H	29	THR	SER	conflict	UNP P02281

- Molecule 5 is a DNA chain called DNA (122-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
5	I	122	2483	1178	451	732	122	0	0

- Molecule 6 is a DNA chain called DNA (122-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
6	J	122	2519	1190	475	732	122	0	0

- Molecule 7 is a protein called TATA-binding protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	K	180	1416	921	242	247	6	0	0

- Molecule 8 is a protein called Transcription initiation factor IIA large subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	L	108	884	550	147	184	3	0	0

There are 115 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	GLY	deletion	UNP A0A6A5Q2T8
L	?	-	LEU	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8

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Chain	Residue	Modelled	Actual	Comment	Reference
L	?	-	LEU	deletion	UNP A0A6A5Q2T8
L	?	-	PRO	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	PRO	deletion	UNP A0A6A5Q2T8
L	?	-	HIS	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	GLY	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	ALA	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	GLY	deletion	UNP A0A6A5Q2T8
L	?	-	ALA	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	LEU	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	GLY	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8

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Chain	Residue	Modelled	Actual	Comment	Reference
L	?	-	ALA	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	GLN	deletion	UNP A0A6A5Q2T8
L	?	-	PRO	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	PRO	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LEU	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ALA	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	ASN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	SER	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8

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Chain	Residue	Modelled	Actual	Comment	Reference
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	ASP	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	THR	deletion	UNP A0A6A5Q2T8
L	?	-	ARG	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	GLN	deletion	UNP A0A6A5Q2T8
L	?	-	ILE	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	GLN	deletion	UNP A0A6A5Q2T8
L	?	-	VAL	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	LEU	deletion	UNP A0A6A5Q2T8
L	?	-	GLN	deletion	UNP A0A6A5Q2T8
L	?	-	ALA	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	GLU	deletion	UNP A0A6A5Q2T8
L	?	-	LYS	deletion	UNP A0A6A5Q2T8
L	?	-	ARG	deletion	UNP A0A6A5Q2T8

- Molecule 9 is a protein called Transcription initiation factor IIA subunit 2.

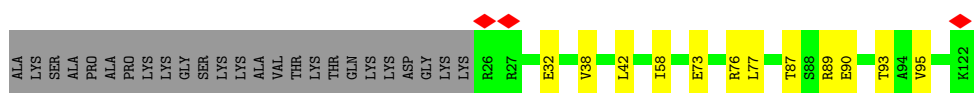
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	M	107	836	522	141	169	4	0	0



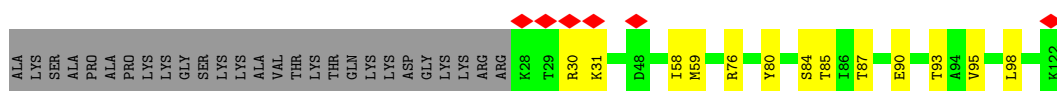
• Molecule 3: Histone H2A



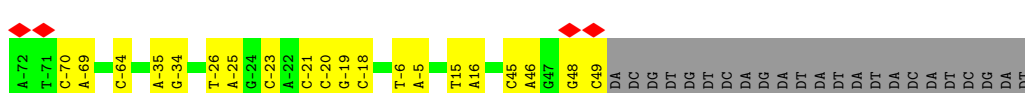
• Molecule 4: Histone H2B 1.1



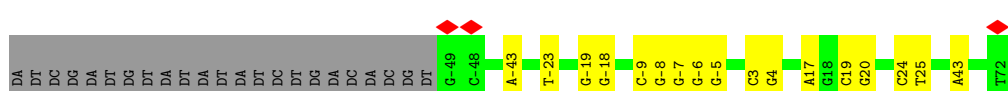
• Molecule 4: Histone H2B 1.1



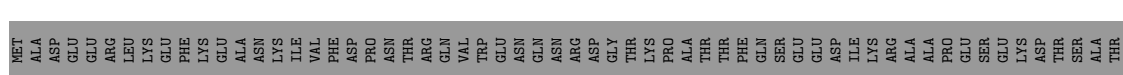
• Molecule 5: DNA (122-MER)

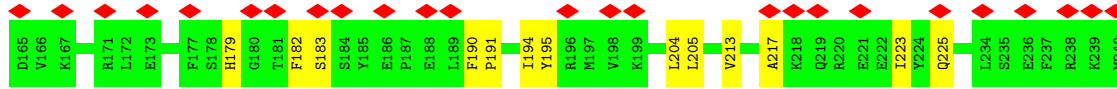


• Molecule 6: DNA (122-MER)

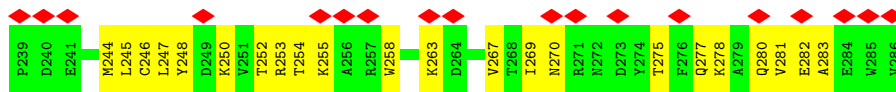
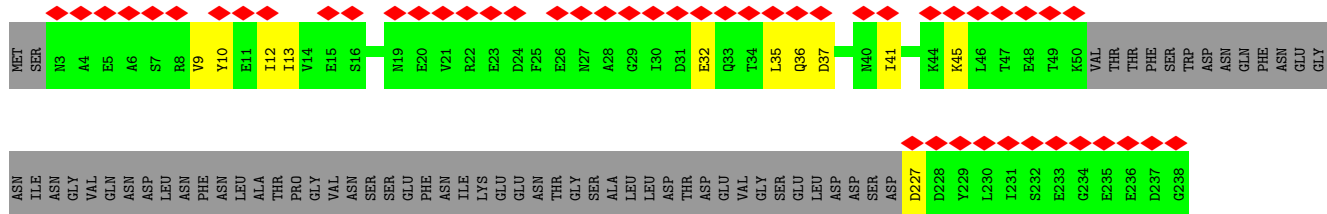
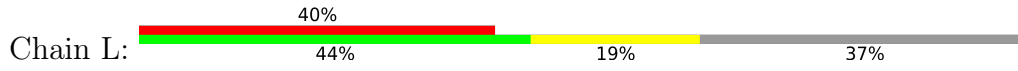


• Molecule 7: TATA-binding protein

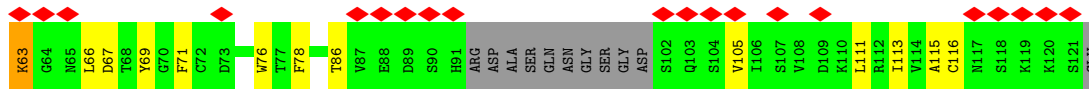
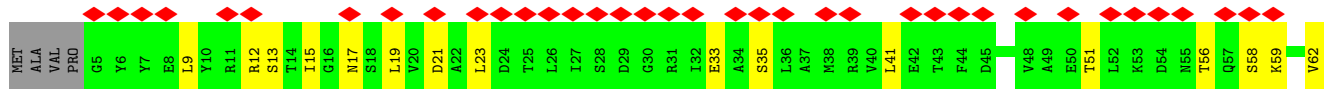




• Molecule 8: Transcription initiation factor IIA large subunit



• Molecule 9: Transcription initiation factor IIA subunit 2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	130350	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 ($e^-/\text{\AA}^2$)	41.2	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.087	Depositor
Minimum map value	-0.032	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.015	Depositor
Map size (\AA)	209.99998, 209.99998, 209.99998	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.05, 1.05, 1.05	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.20	0/814	0.46	0/1092
1	E	0.18	0/823	0.41	0/1104
2	B	0.19	0/660	0.46	0/883
2	F	0.21	0/645	0.53	0/862
3	C	0.16	0/853	0.38	0/1149
3	G	0.17	0/828	0.42	0/1117
4	D	0.17	0/778	0.41	0/1043
4	H	0.21	0/756	0.51	0/1015
5	I	0.21	0/2781	0.50	0/4285
6	J	0.21	0/2829	0.46	0/4369
7	K	0.25	0/1443	0.63	2/1942 (0.1%)
8	L	0.37	0/895	0.82	0/1207
9	M	0.39	0/844	0.94	1/1139 (0.1%)
All	All	0.23	0/14949	0.54	3/21207 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
7	K	225	GLN	N-CA-CB	7.25	120.89	110.16
9	M	63	LYS	CA-CB-CG	5.54	125.19	114.10
7	K	225	GLN	CA-CB-CG	5.06	124.22	114.10

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	802	0	841	12	0
1	E	811	0	853	9	0
2	B	653	0	696	13	0
2	F	638	0	676	5	0
3	C	843	0	908	9	0
3	G	818	0	877	11	0
4	D	767	0	799	11	0
4	H	745	0	773	11	0
5	I	2483	0	1368	15	0
6	J	2519	0	1368	12	0
7	K	1416	0	1493	21	0
8	L	884	0	851	34	0
9	M	836	0	837	24	0
All	All	14215	0	12340	133	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 5.

All (133) 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)
8:L:283:ALA:HB1	9:M:66:LEU:HB2	1.63	0.78
9:M:78:PHE:HB2	9:M:113:ILE:HB	1.67	0.77
8:L:278:LYS:HE3	9:M:59:LYS:HZ3	1.57	0.70
3:G:16:THR:HA	6:J:-43:DA:H5''	1.76	0.68
8:L:9:VAL:HG11	8:L:275:THR:HG21	1.76	0.67
7:K:156:LYS:HE2	7:K:158:GLN:HE21	1.59	0.66
8:L:258:TRP:HB3	8:L:283:ALA:HB3	1.75	0.66
8:L:13:ILE:HG13	9:M:51:THR:HG21	1.78	0.66
2:F:92:ARG:HH21	4:H:98:LEU:HD23	1.63	0.64
3:G:42:ARG:NH1	4:H:85:THR:OG1	2.32	0.63
8:L:32:GLU:HA	8:L:35:LEU:HD23	1.83	0.60
1:A:63:ARG:HD3	6:J:17:DA:H4'	1.84	0.60
7:K:195:TYR:HB3	7:K:204:LEU:HB2	1.83	0.60
4:H:30:ARG:HE	4:H:31:LYS:H	1.50	0.59
7:K:217:ALA:HB2	7:K:223:ILE:HG23	1.84	0.59
1:A:83:ARG:HB2	2:B:80:THR:HG22	1.85	0.58
5:I:-19:DG:OP1	7:K:110:LYS:NZ	2.37	0.57
7:K:205:LEU:HB2	7:K:213:VAL:HB	1.86	0.57
7:K:95:ASN:HB2	9:M:71:PHE:HE2	1.70	0.56
4:D:73:GLU:OE2	4:D:76:ARG:NH2	2.39	0.56
3:G:85:LEU:O	3:G:89:ASN:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:9:VAL:O	8:L:13:ILE:HD12	2.06	0.56
2:B:75:HIS:O	4:D:89:ARG:NH2	2.39	0.55
5:I:48:DG:H1'	5:I:49:DC:H5'	1.89	0.54
3:C:63:LEU:HD13	4:D:42:LEU:HB2	1.90	0.54
2:B:47:SER:OG	2:B:48:GLY:N	2.41	0.54
7:K:107:ARG:NH2	9:M:67:ASP:O	2.41	0.54
3:C:102:ILE:HG23	4:D:58:ILE:HD13	1.89	0.53
7:K:145:LYS:HE2	8:L:227:ASP:HB2	1.90	0.53
5:I:-20:DC:OP1	8:L:253:ARG:NH2	2.42	0.53
8:L:275:THR:O	9:M:58:SER:OG	2.25	0.53
2:B:22:LEU:HD21	8:L:250:LYS:HE3	1.91	0.53
8:L:9:VAL:O	8:L:12:ILE:HG22	2.10	0.52
3:C:63:LEU:HD11	4:D:38:VAL:HG13	1.92	0.52
9:M:19:LEU:O	9:M:23:LEU:HD12	2.10	0.51
7:K:183:SER:OG	7:K:194:ILE:O	2.27	0.51
1:A:128:ARG:HD3	1:A:134:ARG:HH21	1.76	0.51
7:K:105:ARG:NH1	9:M:69:TYR:OH	2.41	0.51
3:C:104:GLN:NE2	1:E:94:GLU:OE1	2.38	0.51
8:L:277:GLN:HB3	9:M:56:THR:HG23	1.93	0.51
2:B:68:ASP:OD2	2:B:92:ARG:NH2	2.45	0.50
8:L:244:MET:HE2	8:L:267:VAL:HG11	1.93	0.50
3:C:113:SER:HA	3:C:116:LEU:HD23	1.94	0.50
3:G:54:VAL:HG21	4:H:95:VAL:HG21	1.94	0.49
5:I:15:DT:H2''	5:I:16:DA:C8	2.48	0.49
5:I:-21:DC:H2'	5:I:-20:DC:C6	2.47	0.49
8:L:245:LEU:HD11	9:M:12:ARG:HB2	1.94	0.49
3:G:83:LEU:HD22	4:H:59:MET:HE3	1.95	0.48
5:I:-18:DC:OP1	7:K:68:GLN:NE2	2.34	0.48
7:K:179:HIS:HB3	7:K:182:PHE:HB3	1.94	0.48
1:E:56:LYS:NZ	5:I:-64:DC:OP1	2.47	0.48
8:L:282:GLU:OE2	9:M:63:LYS:HD2	2.14	0.47
6:J:24:DC:H2''	6:J:25:DT:C6	2.50	0.47
2:B:38:ALA:HB1	2:B:43:VAL:HB	1.95	0.47
3:G:16:THR:HG23	3:G:19:SER:H	1.78	0.47
2:F:75:HIS:HB2	4:H:93:THR:HG21	1.95	0.47
2:B:75:HIS:HB2	4:D:93:THR:HG21	1.95	0.47
6:J:19:DC:H2'	6:J:20:DG:C8	2.50	0.47
7:K:70:ILE:HG23	7:K:160:ILE:HG22	1.97	0.47
6:J:19:DC:H2''	6:J:20:DG:H5'	1.96	0.46
7:K:67:LEU:HD11	7:K:160:ILE:HD12	1.97	0.46
1:A:46:VAL:O	1:A:50:GLU:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:247:LEU:O	8:L:263:LYS:N	2.36	0.46
4:H:76:ARG:HG2	4:H:80:TYR:CZ	2.50	0.46
4:H:87:THR:OG1	4:H:90:GLU:OE1	2.34	0.46
8:L:258:TRP:N	8:L:283:ALA:O	2.49	0.46
5:I:-35:DA:H2''	5:I:-34:DG:C8	2.50	0.46
5:I:-26:DT:H2''	5:I:-25:DA:C8	2.49	0.46
9:M:86:THR:HG22	9:M:105:VAL:HG22	1.97	0.46
6:J:3:DC:H2''	6:J:4:DG:C8	2.50	0.45
1:A:61:LEU:HD12	2:B:37:LEU:HD23	1.98	0.45
8:L:32:GLU:O	8:L:36:GLN:OE1	2.34	0.45
8:L:269:ILE:HG22	8:L:270:ASN:OD1	2.16	0.45
7:K:190:PHE:CD1	7:K:191:PRO:HD2	2.52	0.45
3:G:102:ILE:HG23	4:H:58:ILE:HD13	1.98	0.45
8:L:244:MET:HG3	8:L:267:VAL:HG12	1.98	0.45
9:M:13:SER:O	9:M:17:ASN:N	2.43	0.45
5:I:45:DC:H2''	5:I:46:DA:C8	2.53	0.44
3:G:68:ASN:OD1	3:G:71:ARG:NH2	2.48	0.44
6:J:-7:DG:H2''	6:J:-6:DG:C8	2.52	0.44
8:L:37:ASP:O	8:L:41:ILE:HG12	2.17	0.44
8:L:253:ARG:NH1	8:L:254:THR:O	2.50	0.44
1:A:113:HIS:CG	1:E:126:LEU:HD22	2.52	0.44
1:A:126:LEU:HD22	1:E:113:HIS:CG	2.53	0.44
6:J:-9:DC:H2''	6:J:-8:DG:C8	2.53	0.44
6:J:-6:DG:H2''	6:J:-5:DG:C8	2.53	0.44
9:M:41:LEU:HD23	9:M:41:LEU:HA	1.79	0.43
1:A:51:ILE:HG13	2:B:39:ARG:HD2	2.01	0.43
3:C:54:VAL:HG21	4:D:95:VAL:HG21	2.00	0.43
2:F:68:ASP:OD2	2:F:92:ARG:NH1	2.52	0.43
3:G:88:ARG:NH2	3:G:100:VAL:O	2.51	0.43
3:C:32:ARG:HH22	4:D:32:GLU:CD	2.27	0.43
9:M:71:PHE:HB2	9:M:76:TRP:CZ3	2.54	0.43
1:A:133:GLU:OE2	2:B:95:ARG:NH2	2.51	0.43
8:L:250:LYS:NZ	8:L:252:THR:OG1	2.40	0.43
2:B:24:ASP:OD1	2:B:25:ASN:N	2.52	0.43
5:I:-6:DT:H2''	5:I:-5:DA:N7	2.34	0.43
3:C:11:ARG:HH12	6:J:43:DA:H2	1.67	0.42
1:E:72:ARG:O	1:E:76:GLN:HG3	2.19	0.42
8:L:45:LYS:HB3	8:L:45:LYS:HE2	1.78	0.42
7:K:130:ASP:HA	7:K:133:LYS:HG2	2.00	0.42
1:A:123:ASP:OD1	1:E:113:HIS:NE2	2.46	0.42
8:L:45:LYS:NZ	9:M:21:ASP:OD1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:250:LYS:HZ2	8:L:252:THR:HG1	1.62	0.42
3:C:85:LEU:O	3:C:89:ASN:HB2	2.20	0.42
5:I:-70:DC:H2''	5:I:-69:DA:C8	2.55	0.42
8:L:248:TYR:CE1	9:M:76:TRP:HD1	2.37	0.42
2:F:75:HIS:NE2	4:H:90:GLU:HG3	2.35	0.42
7:K:133:LYS:O	7:K:137:ARG:HG2	2.19	0.42
8:L:281:VAL:HG12	9:M:62:VAL:HB	2.02	0.42
6:J:-19:DG:H2''	6:J:-18:DG:C8	2.55	0.42
1:A:42:ARG:NH1	5:I:-5:DA:OP1	2.53	0.42
1:A:72:ARG:NH1	5:I:-23:DC:OP1	2.50	0.42
5:I:-18:DC:H41	8:L:255:LYS:NZ	2.18	0.42
4:D:77:LEU:HD21	4:D:93:THR:HB	2.01	0.41
1:E:46:VAL:O	1:E:50:GLU:HG3	2.20	0.41
7:K:82:LEU:HD11	7:K:117:ALA:HA	2.02	0.41
7:K:130:ASP:OD1	7:K:131:ASP:N	2.53	0.41
8:L:244:MET:HE3	9:M:111:LEU:HD21	2.02	0.41
2:B:22:LEU:HD22	8:L:280:GLN:HE22	1.85	0.41
9:M:9:LEU:HD12	9:M:116:CYS:SG	2.61	0.41
2:F:31:LYS:HE2	2:F:51:TYR:CE2	2.56	0.41
7:K:70:ILE:HG12	7:K:126:ALA:O	2.20	0.41
1:E:83:ARG:HG2	6:J:-23:DT:H5''	2.02	0.41
9:M:33:GLU:CD	9:M:35:SER:H	2.29	0.41
7:K:142:ILE:O	7:K:146:ILE:HG12	2.21	0.41
8:L:246:CYS:O	9:M:115:ALA:HA	2.21	0.41
2:B:75:HIS:CE1	4:D:77:LEU:HG	2.56	0.41
8:L:10:TYR:CZ	9:M:15:ILE:HD11	2.56	0.40
4:D:87:THR:OG1	4:D:90:GLU:OE1	2.29	0.40
3:G:41:GLU:HG3	4:H:84:SER:HB2	2.04	0.40
1:E:128:ARG:HB2	1:E:134:ARG:HD2	2.02	0.40
3:G:41:GLU:OE1	3:G:42:ARG:HG3	2.22	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	95/135 (70%)	94 (99%)	1 (1%)	0	100	100
1	E	96/135 (71%)	93 (97%)	3 (3%)	0	100	100
2	B	80/102 (78%)	78 (98%)	2 (2%)	0	100	100
2	F	78/102 (76%)	76 (97%)	2 (3%)	0	100	100
3	C	107/129 (83%)	106 (99%)	1 (1%)	0	100	100
3	G	104/129 (81%)	102 (98%)	2 (2%)	0	100	100
4	D	95/122 (78%)	93 (98%)	2 (2%)	0	100	100
4	H	93/122 (76%)	93 (100%)	0	0	100	100
7	K	178/240 (74%)	177 (99%)	1 (1%)	0	100	100
8	L	104/171 (61%)	102 (98%)	2 (2%)	0	100	100
9	M	103/122 (84%)	103 (100%)	0	0	100	100
All	All	1133/1509 (75%)	1117 (99%)	16 (1%)	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	85/110 (77%)	85 (100%)	0	100	100
1	E	86/110 (78%)	86 (100%)	0	100	100
2	B	67/78 (86%)	67 (100%)	0	100	100
2	F	65/78 (83%)	65 (100%)	0	100	100
3	C	86/101 (85%)	86 (100%)	0	100	100
3	G	84/101 (83%)	84 (100%)	0	100	100
4	D	83/102 (81%)	83 (100%)	0	100	100
4	H	81/102 (79%)	81 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	K	152/205 (74%)	152 (100%)	0	100	100
8	L	98/154 (64%)	98 (100%)	0	100	100
9	M	97/108 (90%)	97 (100%)	0	100	100
All	All	984/1249 (79%)	984 (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 (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
3	C	73	ASN
4	D	64	ASN
4	D	79	HIS
4	D	106	HIS
1	E	68	GLN
1	E	108	ASN
3	G	24	GLN
3	G	38	ASN
3	G	73	ASN
4	H	46	HIS
7	K	158	GLN
8	L	280	GLN
9	M	57	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

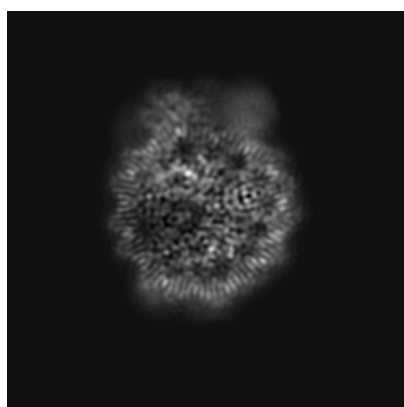
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12898. These allow visual inspection of the internal detail of the map and identification of artifacts.

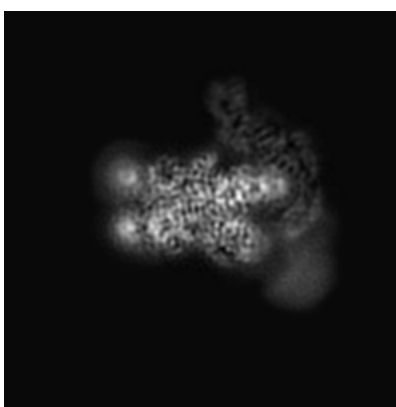
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

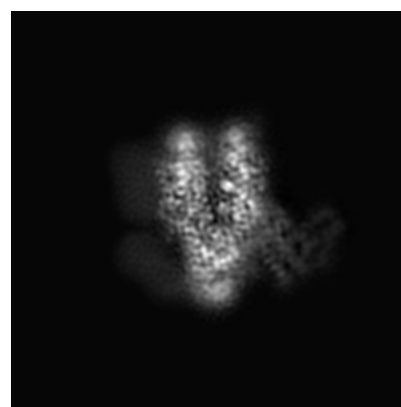
6.1.1 Primary map



X



Y

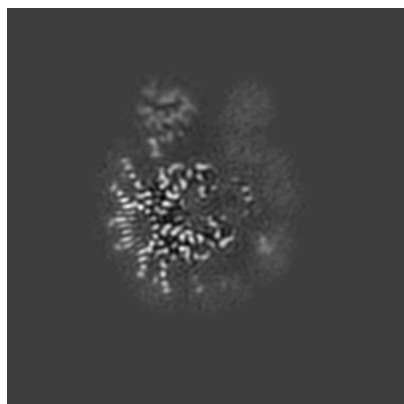


Z

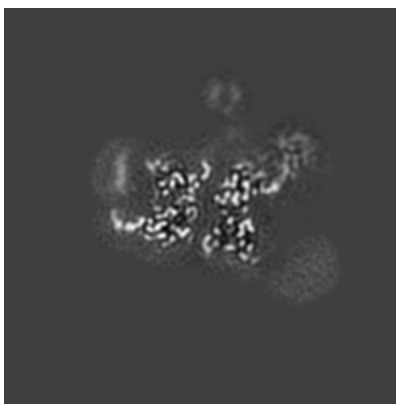
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

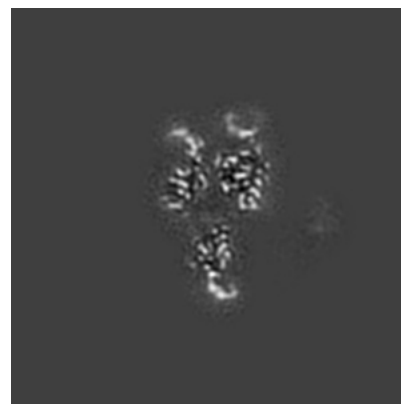
6.2.1 Primary map



X Index: 100



Y Index: 100

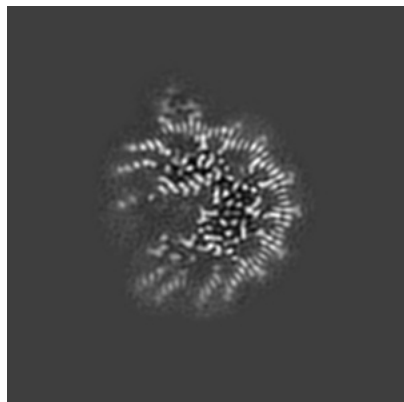


Z Index: 100

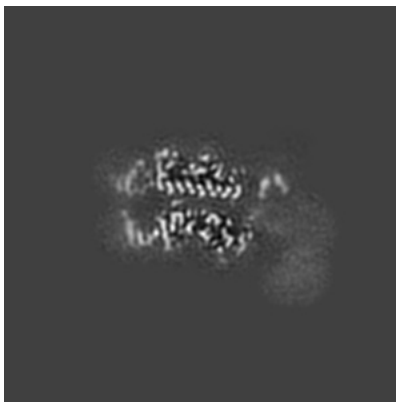
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

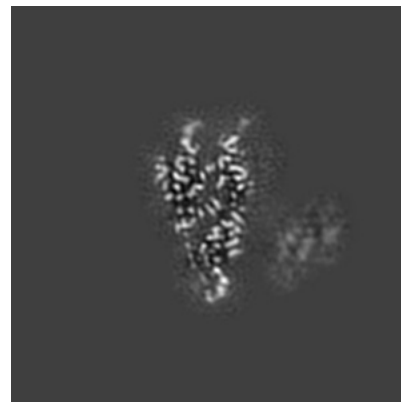
6.3.1 Primary map



X Index: 112



Y Index: 112

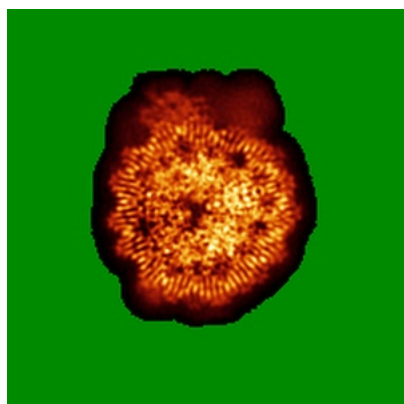


Z Index: 109

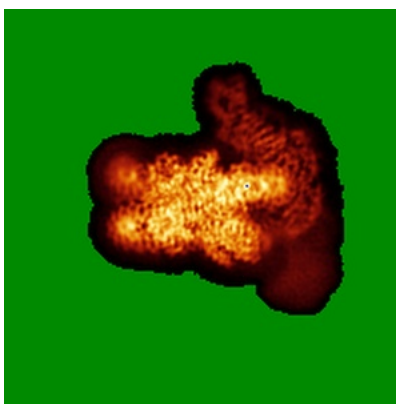
The images above show the largest variance slices of the map in three orthogonal directions.

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

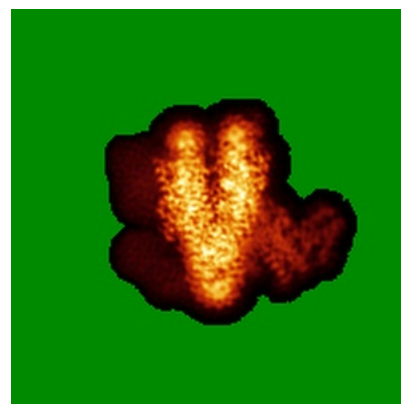
6.4.1 Primary map



X



Y

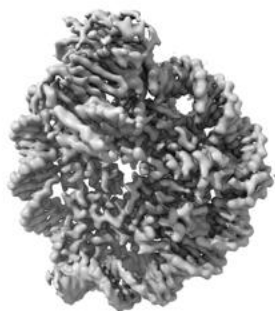


Z

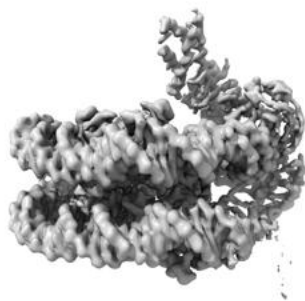
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

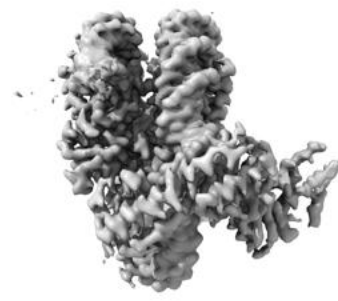
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.015. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

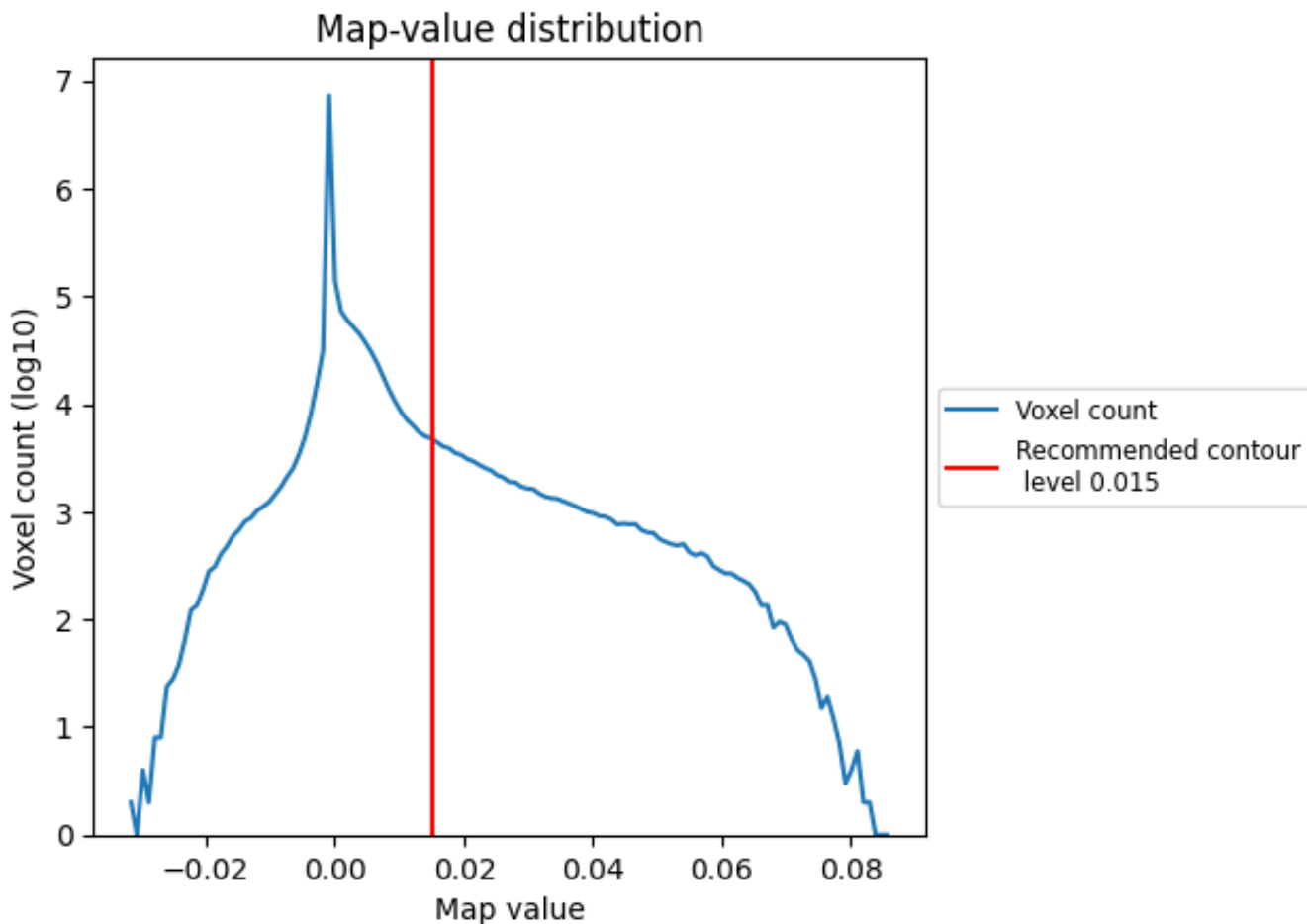
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

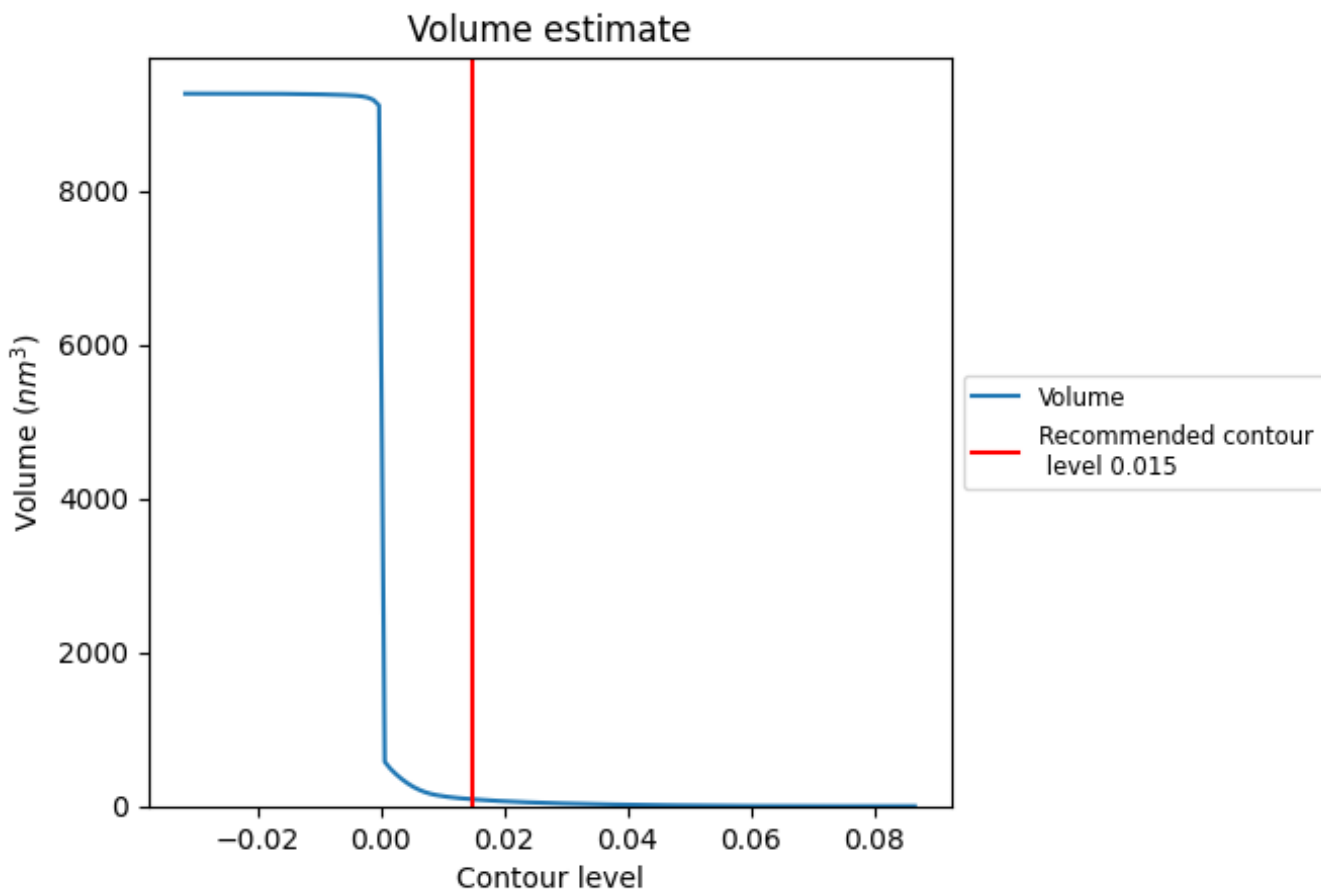
This section contains the results of statistical analysis of the map.

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.

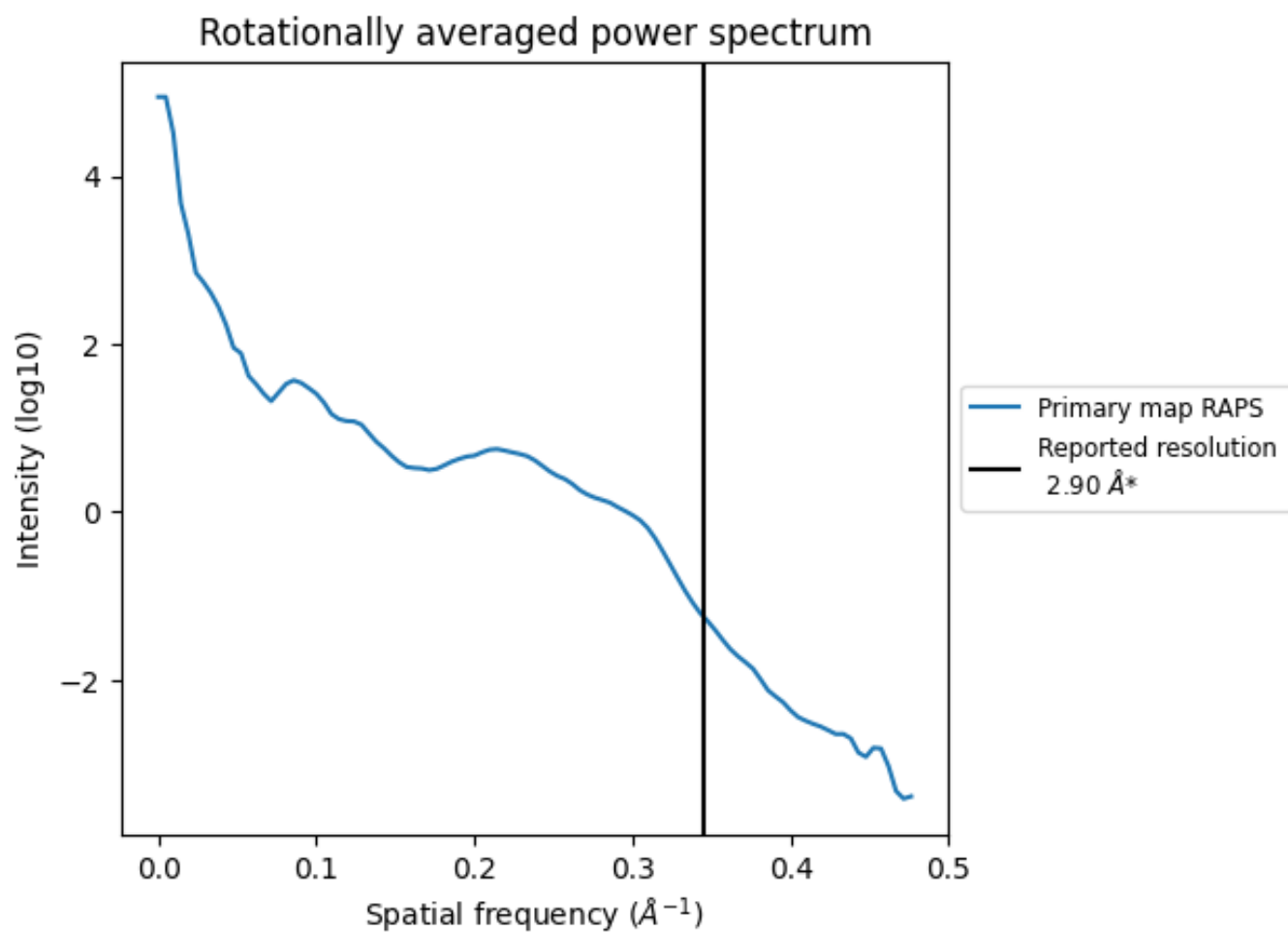
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 89 nm^3 ; this corresponds to an approximate mass of 81 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

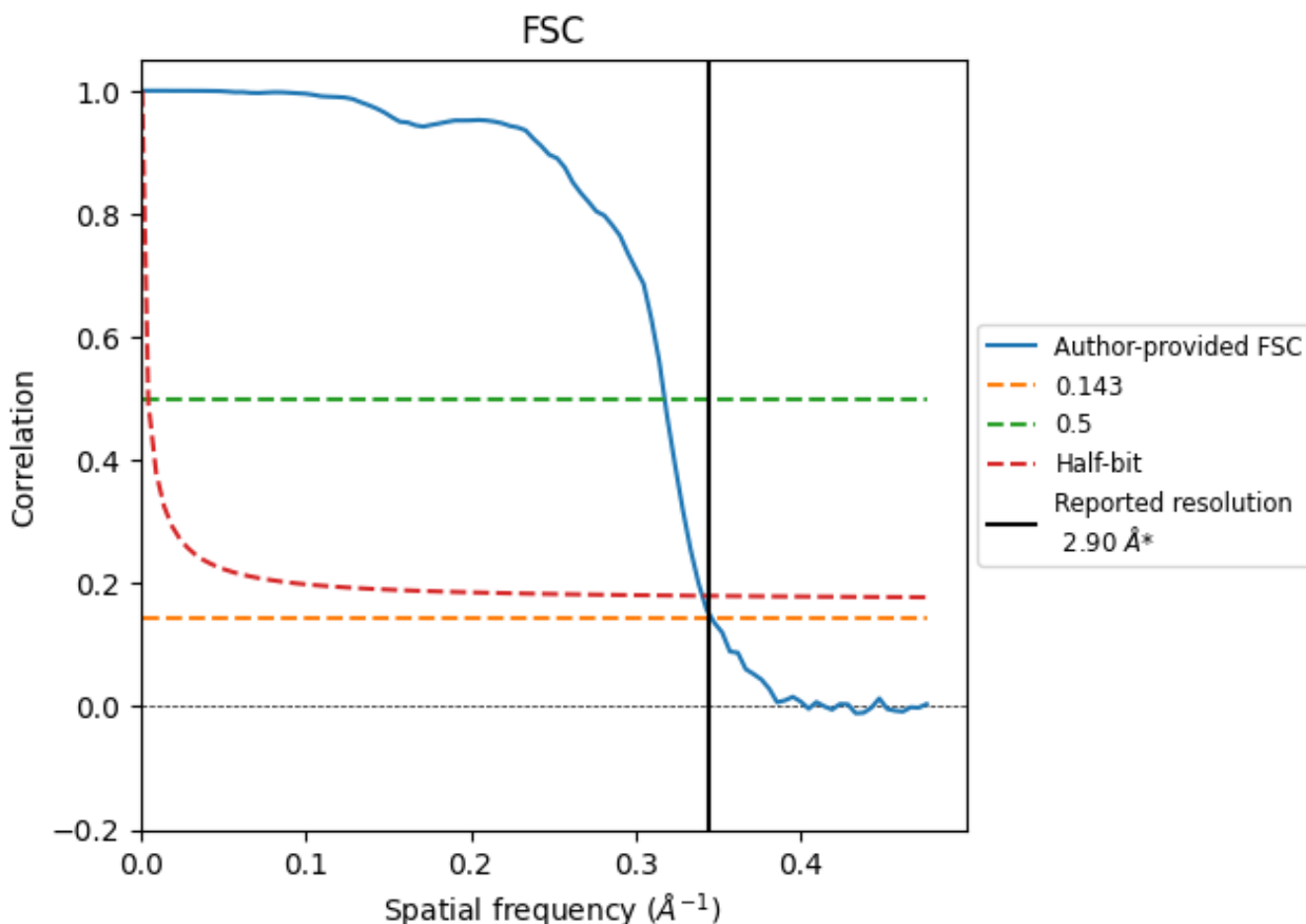


*Reported resolution corresponds to spatial frequency of 0.345 Å⁻¹

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 0.345 Å⁻¹

8.2 Resolution estimates [i](#)

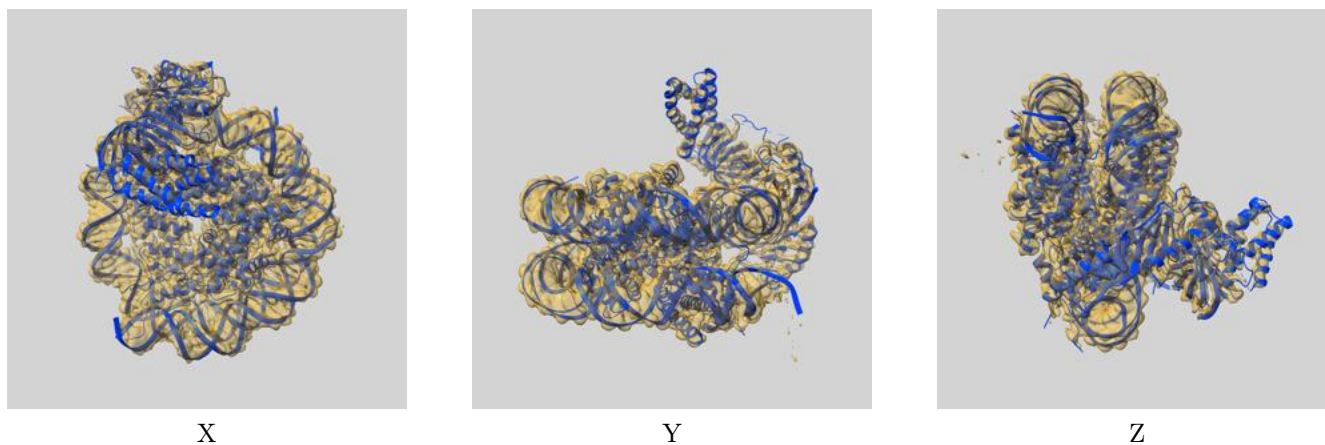
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	2.89	3.15	2.94
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

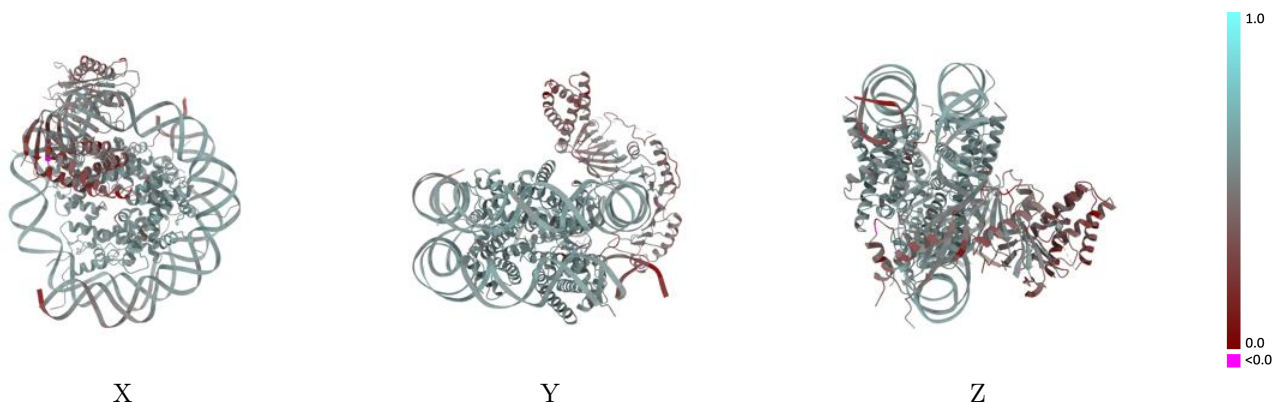
This section contains information regarding the fit between EMDB map EMD-12898 and PDB model 7OHA. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



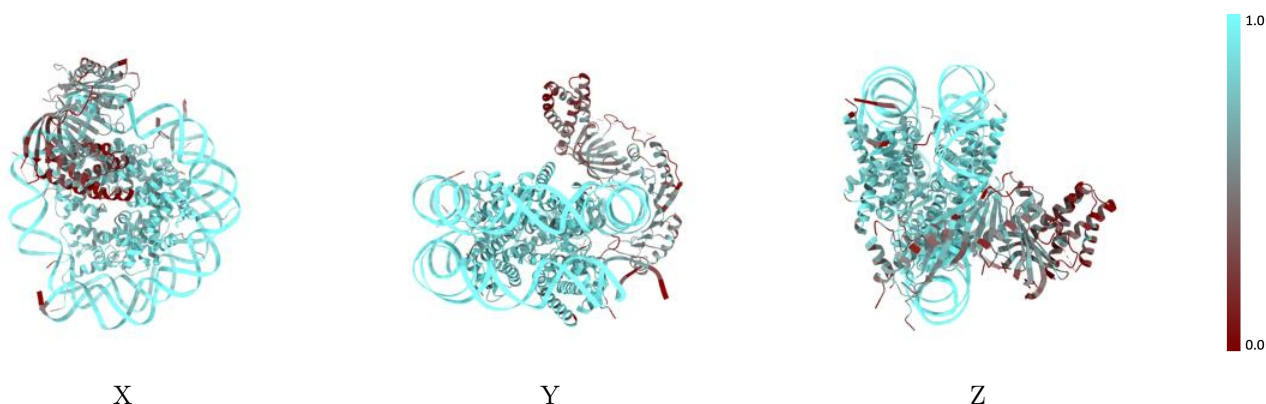
The images above show the 3D surface view of the map at the recommended contour level 0.015 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



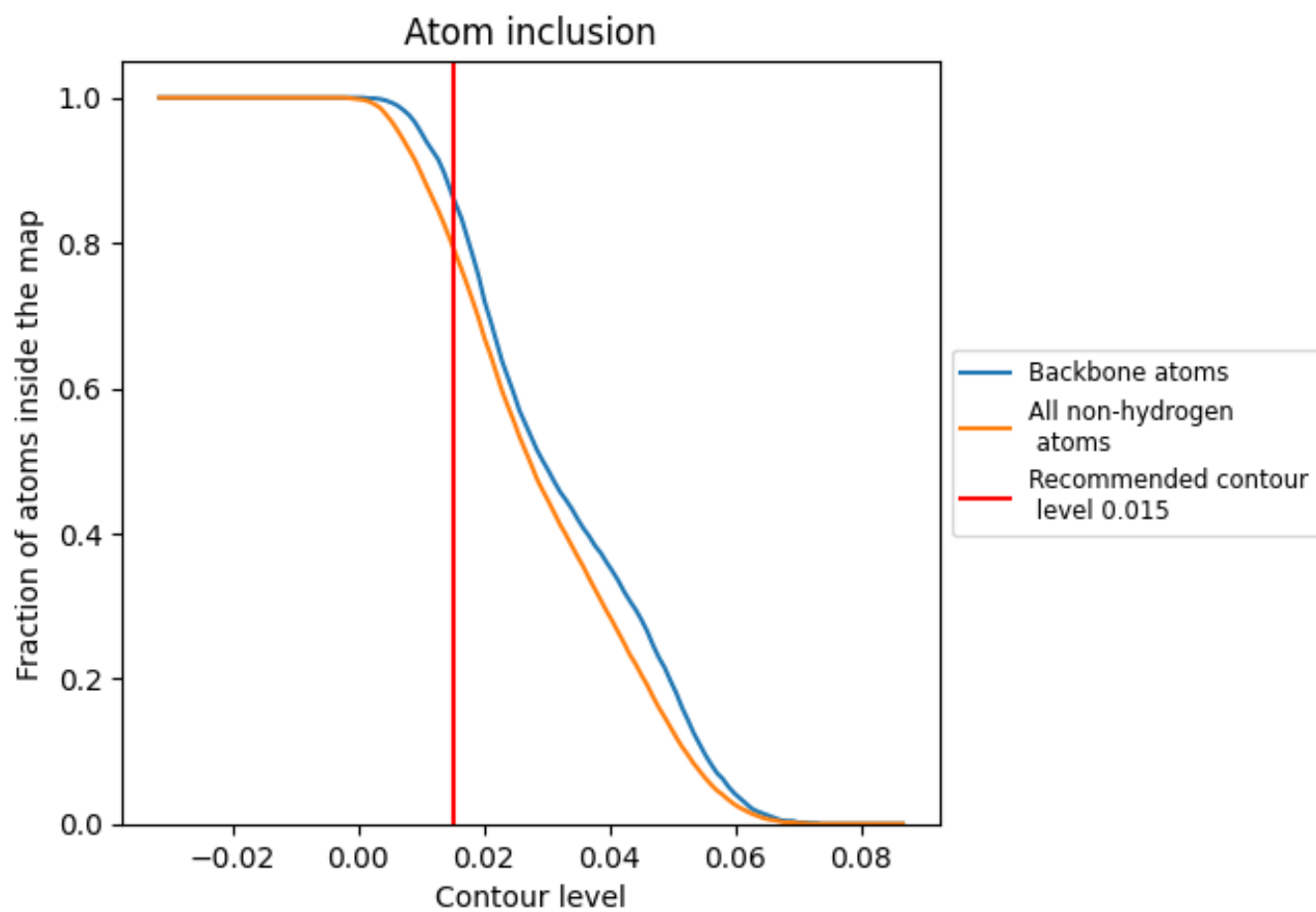
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.015).





























9.4 Atom inclusion [i](#)



At the recommended contour level, 86% 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 (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7960	 0.5160
A	 0.8530	 0.5470
B	 0.8660	 0.5580
C	 0.8930	 0.5740
D	 0.8810	 0.5540
E	 0.9050	 0.5760
F	 0.9260	 0.5890
G	 0.8500	 0.5650
H	 0.8310	 0.5340
I	 0.9370	 0.5440
J	 0.9340	 0.5420
K	 0.5130	 0.4200
L	 0.3020	 0.3390
M	 0.3740	 0.3660

