



wwPDB EM Validation Summary Report ⓘ

Mar 10, 2026 – 05:23 AM UTC

PDB ID : 8OTS / pdb_00008ots
EMDB ID : EMD-17183
Title : OCT4 and MYC-MAX co-bound to a nucleosome
Authors : Michael, A.K.; Stoos, L.; Kempf, G.; Cavadini, S.; Thoma, N.
Deposited on : 2023-04-21
Resolution : 3.30 Å(reported)
Based on initial models : 6T90, 1NKP

This is a wwPDB EM Validation Summary 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
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

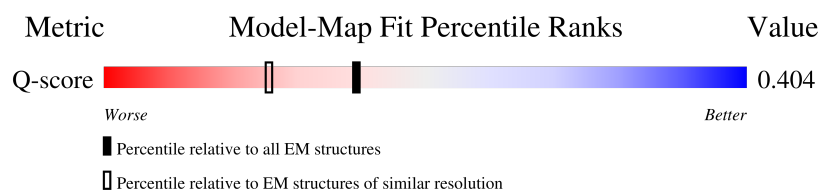
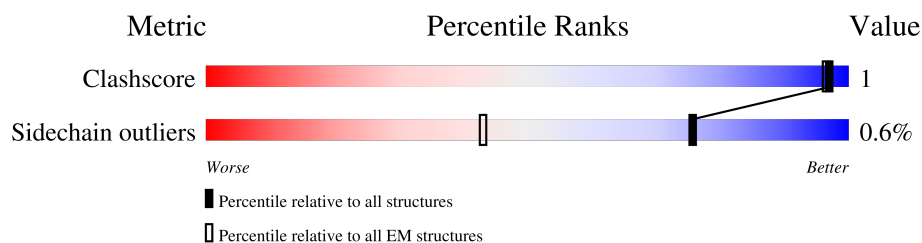
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15087 (2.80 - 3.80)

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	139	
1	E	139	
2	B	106	
2	F	106	
3	C	133	

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Mol	Chain	Length	Quality of chain
3	G	133	
4	D	128	
4	H	128	
5	I	127	
6	J	127	
7	K	544	
8	M	94	
9	N	82	

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13160 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 H3.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	94	Total	C	N	O	S	0	0
			773	488	147	134	4		
1	E	95	Total	C	N	O	S	0	0
			783	494	150	135	4		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P68431
A	-2	SER	-	expression tag	UNP P68431
A	-1	HIS	-	expression tag	UNP P68431
E	-3	GLY	-	expression tag	UNP P68431
E	-2	SER	-	expression tag	UNP P68431
E	-1	HIS	-	expression tag	UNP P68431

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	82	Total	C	N	O	S	0	0
			653	412	127	113	1		
2	F	81	Total	C	N	O	S	0	0
			646	407	126	112	1		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P62805
B	-2	SER	-	expression tag	UNP P62805
B	-1	HIS	-	expression tag	UNP P62805
F	-3	GLY	-	expression tag	UNP P62805
F	-2	SER	-	expression tag	UNP P62805
F	-1	HIS	-	expression tag	UNP P62805

- Molecule 3 is a protein called Histone H2A type 1-B/E.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	109	Total	C	N	O	0	0
			837	526	166	145		
3	G	109	Total	C	N	O	0	0
			840	529	166	145		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLY	-	expression tag	UNP P04908
C	-2	SER	-	expression tag	UNP P04908
C	-1	HIS	-	expression tag	UNP P04908
G	-3	GLY	-	expression tag	UNP P04908
G	-2	SER	-	expression tag	UNP P04908
G	-1	HIS	-	expression tag	UNP P04908

- Molecule 4 is a protein called Histone H2B type 1-J.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	93	Total	C	N	O	S	0	0
			725	456	130	137	2		
4	H	92	Total	C	N	O	S	0	0
			719	453	129	135	2		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-3	GLY	-	expression tag	UNP P06899
D	-2	SER	-	expression tag	UNP P06899
D	-1	HIS	-	expression tag	UNP P06899
H	-3	GLY	-	expression tag	UNP P06899
H	-2	SER	-	expression tag	UNP P06899
H	-1	HIS	-	expression tag	UNP P06899

- Molecule 5 is a DNA chain called DNA (127-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	127	Total	C	N	O	P	0	0
			2592	1232	469	765	126		

- Molecule 6 is a DNA chain called DNA (127-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	127	Total	C	N	O	P	0	0
			2609	1236	492	755	126		

- Molecule 7 is a protein called Green fluorescent protein,POU domain, class 5, transcription factor 1.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	K	73	Total	C	N	O	0	0
			363	217	73	73		

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	-270	MET	-	initiating methionine	UNP P42212
K	-269	ASP	-	expression tag	UNP P42212
K	-268	TRP	-	expression tag	UNP P42212
K	-267	SER	-	expression tag	UNP P42212
K	-266	HIS	-	expression tag	UNP P42212
K	-265	PRO	-	expression tag	UNP P42212
K	-264	GLN	-	expression tag	UNP P42212
K	-263	PHE	-	expression tag	UNP P42212
K	-262	GLU	-	expression tag	UNP P42212
K	-261	LYS	-	expression tag	UNP P42212
K	-260	SER	-	expression tag	UNP P42212
K	-259	ALA	-	expression tag	UNP P42212
K	-258	VAL	-	expression tag	UNP P42212
K	-257	ASP	-	expression tag	UNP P42212
K	-256	GLU	-	expression tag	UNP P42212
K	-255	ASN	-	expression tag	UNP P42212
K	-254	LEU	-	expression tag	UNP P42212
K	-253	TYR	-	expression tag	UNP P42212
K	-252	PHE	-	expression tag	UNP P42212
K	-251	GLN	-	expression tag	UNP P42212
K	-250	GLY	-	expression tag	UNP P42212
K	-249	GLY	-	expression tag	UNP P42212
K	-248	MET	-	expression tag	UNP P42212
K	-247	VAL	-	expression tag	UNP P42212
K	-184	LEU	PHE	conflict	UNP P42212
K	-183	THR	SER	conflict	UNP P42212
K	-42	LYS	ALA	conflict	UNP P42212
K	-17	LEU	HIS	conflict	UNP P42212
K	-9	GLU	-	linker	UNP P42212
K	-8	ALA	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
K	-7	ALA	-	linker	UNP P42212
K	-6	ALA	-	linker	UNP P42212
K	-5	LYS	-	linker	UNP P42212
K	-4	GLU	-	linker	UNP P42212
K	-3	ALA	-	linker	UNP P42212
K	-2	ALA	-	linker	UNP P42212
K	-1	ALA	-	linker	UNP P42212
K	0	LYS	-	linker	UNP P42212

- Molecule 8 is a protein called Myc proto-oncogene protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	M	79	Total	C	N	O	79	0
			798	482	158	158		

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	344	MET	-	initiating methionine	UNP P01106
M	345	HIS	-	expression tag	UNP P01106
M	346	HIS	-	expression tag	UNP P01106
M	347	HIS	-	expression tag	UNP P01106
M	348	HIS	-	expression tag	UNP P01106
M	349	HIS	-	expression tag	UNP P01106
M	350	HIS	-	expression tag	UNP P01106

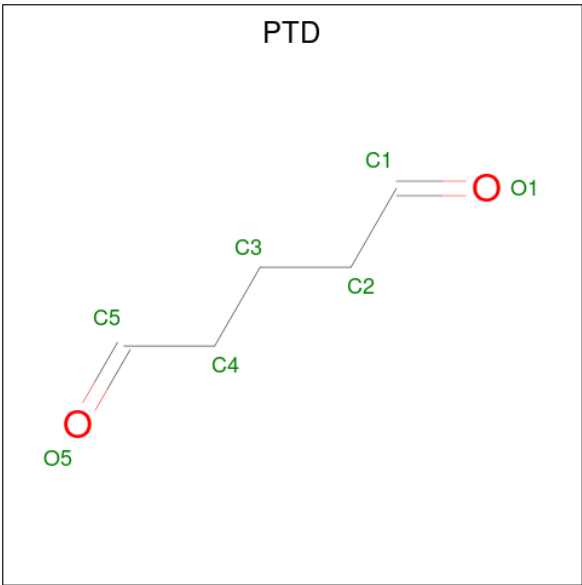
- Molecule 9 is a protein called Protein max.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	N	78	Total	C	N	O	78	0
			782	470	156	156		

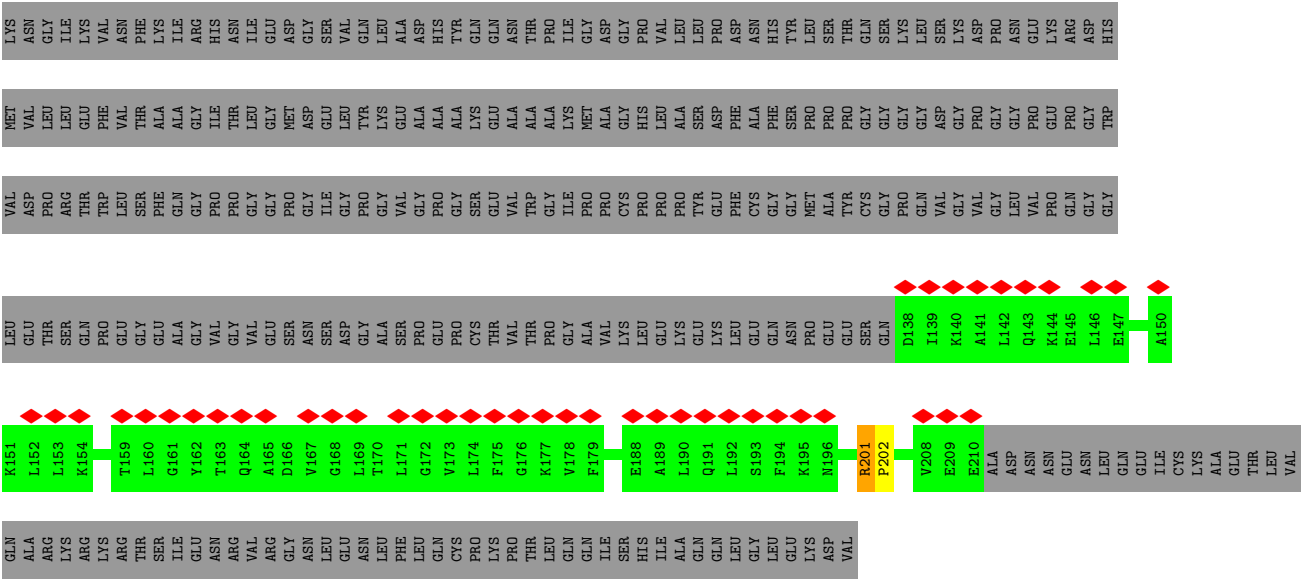
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	21	MET	-	initiating methionine	UNP P61244

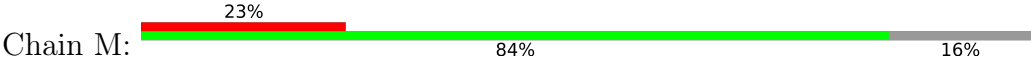
- Molecule 10 is PENTANEDIAL (CCD ID: PTD) (formula: C₅H₈O₂).



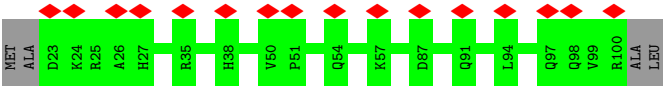
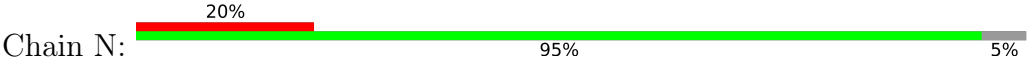
Mol	Chain	Residues	Atoms	AltConf
10	A	1	Total C 5 5	0
10	D	1	Total C 5 5	0
10	E	1	Total C 5 5	0
10	E	1	Total C 5 5	0
10	E	1	Total C 5 5	0
10	G	1	Total C 5 5	0
10	H	1	Total C 5 5	0
10	H	1	Total C 5 5	0



• Molecule 8: Myc proto-oncogene protein



• Molecule 9: Protein max



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	102411	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.667	Depositor
Minimum map value	-0.259	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.025	Depositor
Recommended contour level	0.116	Depositor
Map size ($\text{\AA}$)	220.16, 220.16, 220.16	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.86, 0.86, 0.86	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PTD

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	0.65	0/783	0.76	6/1050 (0.6%)
1	E	0.63	0/794	0.76	4/1065 (0.4%)
2	B	0.61	0/660	0.75	2/883 (0.2%)
2	F	0.62	0/653	0.75	2/873 (0.2%)
3	C	0.61	0/846	0.76	6/1139 (0.5%)
3	G	0.60	0/850	0.77	4/1146 (0.3%)
4	D	0.58	0/736	0.76	4/990 (0.4%)
4	H	0.57	0/730	0.76	4/982 (0.4%)
5	I	0.31	0/2904	0.66	1/4479 (0.0%)
6	J	0.30	0/2930	0.60	0/4522
7	K	0.39	0/363	0.78	2/505 (0.4%)
8	M	0.31	0/800	0.57	0/1120
9	N	0.30	0/782	0.55	0/1092
All	All	0.46	0/13831	0.68	35/19846 (0.2%)

There are no bond length outliers.

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	49	HIS	CA-C-N	7.73	127.44	119.56
4	H	49	HIS	C-N-CA	7.73	127.44	119.56
4	D	49	HIS	CA-C-N	7.30	127.00	119.56
4	D	49	HIS	C-N-CA	7.30	127.00	119.56
4	D	102	LEU	CA-C-N	6.34	126.86	120.14

There are no chirality outliers.

There are no planarity outliers.

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	773	0	804	1	0
1	E	783	0	811	2	0
2	B	653	0	694	2	0
2	F	646	0	685	2	0
3	C	837	0	895	4	0
3	G	840	0	900	3	0
4	D	725	0	739	0	0
4	H	719	0	733	1	0
5	I	2592	0	1429	0	0
6	J	2609	0	1424	1	0
7	K	363	0	166	1	0
8	M	798	0	291	0	0
9	N	782	0	263	0	0
10	A	5	0	6	0	0
10	D	5	0	6	0	0
10	E	15	0	18	0	0
10	G	5	0	6	0	0
10	H	10	0	12	0	0
All	All	13160	0	9882	15	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 1.

The worst 5 of 15 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:E:63:ARG:HH11	6:J:60:DA:H5''	1.68	0.58
3:G:71:ARG:HH22	4:H:49:HIS:CE1	2.27	0.53
3:C:47:ALA:N	3:C:48:PRO:HD2	2.24	0.53
3:G:47:ALA:N	3:G:48:PRO:HD2	2.23	0.52
1:E:65:LEU:N	1:E:66:PRO:HD2	2.28	0.49

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	82/113 (73%)	82 (100%)	0	100	100
1	E	83/113 (74%)	83 (100%)	0	100	100
2	B	67/81 (83%)	67 (100%)	0	100	100
2	F	66/81 (82%)	65 (98%)	1 (2%)	57	72
3	C	84/102 (82%)	83 (99%)	1 (1%)	63	75
3	G	85/102 (83%)	84 (99%)	1 (1%)	63	75
4	D	79/106 (74%)	78 (99%)	1 (1%)	61	74
4	H	78/106 (74%)	78 (100%)	0	100	100
7	K	1/455 (0%)	1 (100%)	0	100	100
8	M	4/89 (4%)	4 (100%)	0	100	100
9	N	2/73 (3%)	2 (100%)	0	100	100
All	All	631/1421 (44%)	627 (99%)	4 (1%)	76	81

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	36	LYS
4	D	85	LYS
2	F	79	LYS
3	G	36	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	24	GLN
3	G	31	HIS

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](#)

8 ligands are 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	$\# Z > 2$
10	PTD	H	202	4	4,4,6	0.44	0	3,3,5	0.58	0
10	PTD	H	201	4,3	4,4,6	0.39	0	3,3,5	0.49	0
10	PTD	E	203	1	4,4,6	0.42	0	3,3,5	0.51	0
10	PTD	E	201	1	4,4,6	0.41	0	3,3,5	0.42	0
10	PTD	A	201	1,2	4,4,6	0.40	0	3,3,5	0.50	0
10	PTD	E	202	1,2	4,4,6	0.38	0	3,3,5	0.51	0
10	PTD	G	200	4,3	4,4,6	0.49	0	3,3,5	0.36	0
10	PTD	D	200	4	4,4,6	0.45	0	3,3,5	0.43	0

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
10	PTD	H	202	4	-	0/2/2/4	-
10	PTD	H	201	4,3	-	0/2/2/4	-
10	PTD	E	203	1	-	0/2/2/4	-
10	PTD	E	201	1	-	1/2/2/4	-
10	PTD	A	201	1,2	-	1/2/2/4	-
10	PTD	E	202	1,2	-	0/2/2/4	-
10	PTD	G	200	4,3	-	2/2/2/4	-
10	PTD	D	200	4	-	2/2/2/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	E	201	PTD	C2-C3-C4-C5
10	D	200	PTD	C1-C2-C3-C4
10	A	201	PTD	C2-C3-C4-C5
10	D	200	PTD	C2-C3-C4-C5
10	G	200	PTD	C2-C3-C4-C5

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.

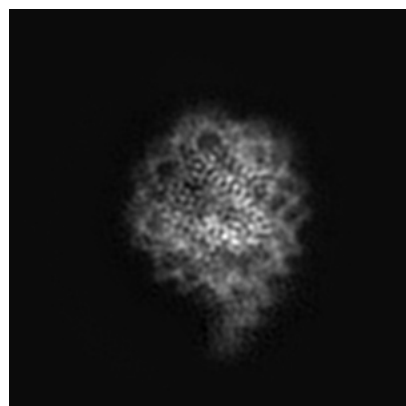
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17183. These allow visual inspection of the internal detail of the map and identification of artifacts.

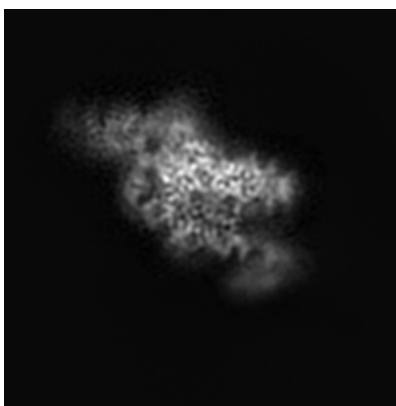
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

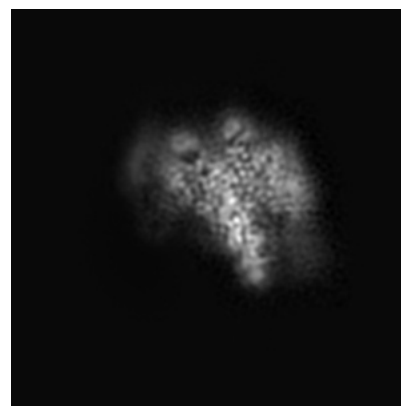
6.1.1 Primary map



X

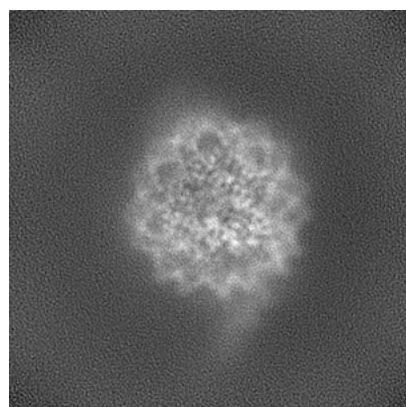


Y

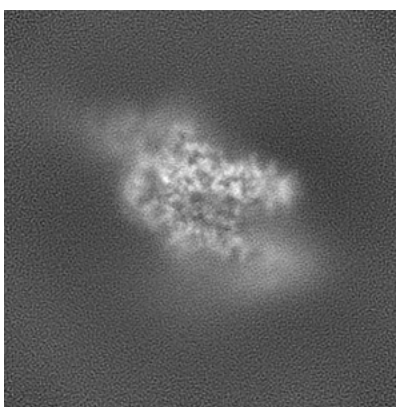


Z

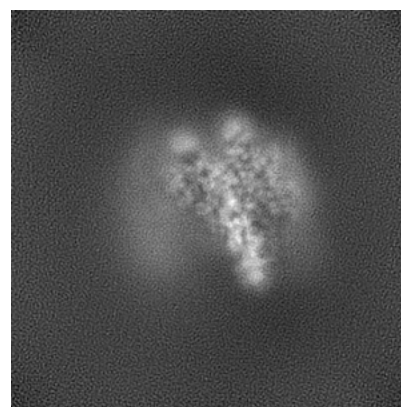
6.1.2 Raw map



X



Y

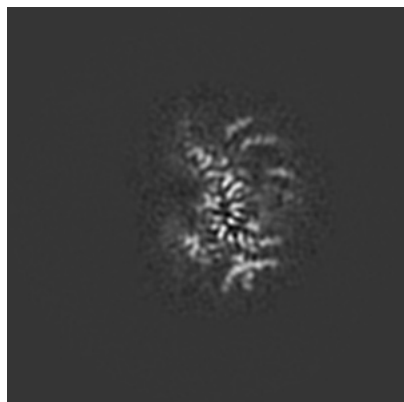


Z

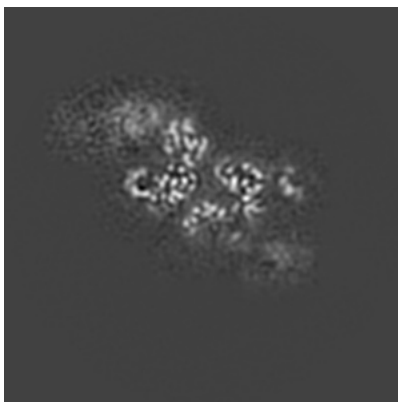
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

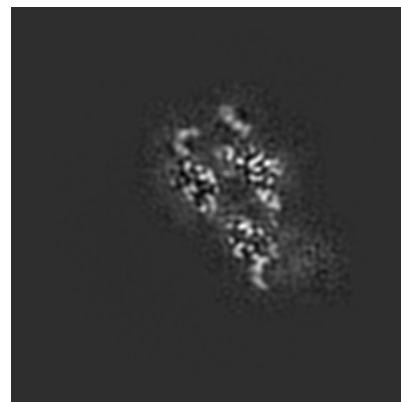
6.2.1 Primary map



X Index: 128

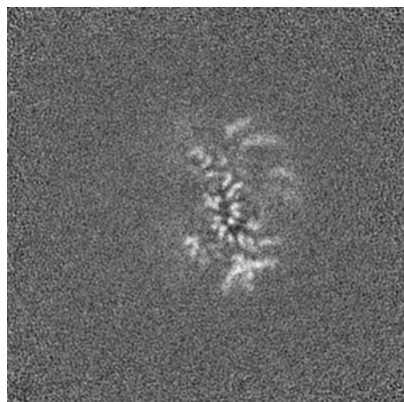


Y Index: 128

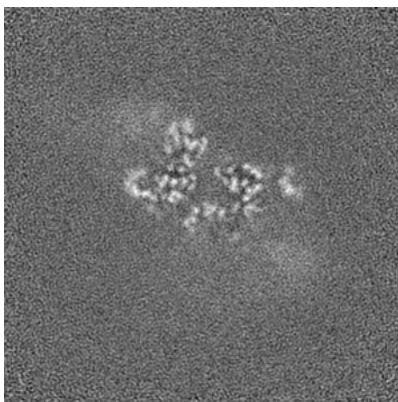


Z Index: 128

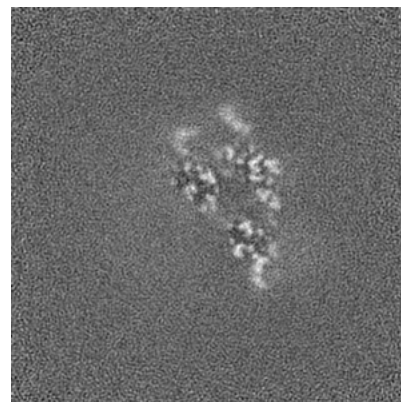
6.2.2 Raw map



X Index: 128



Y Index: 128

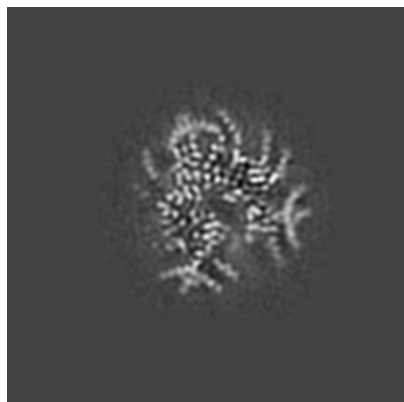


Z Index: 128

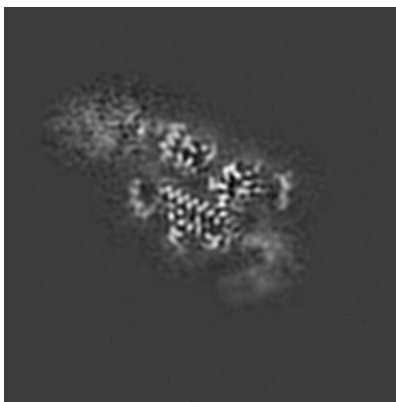
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

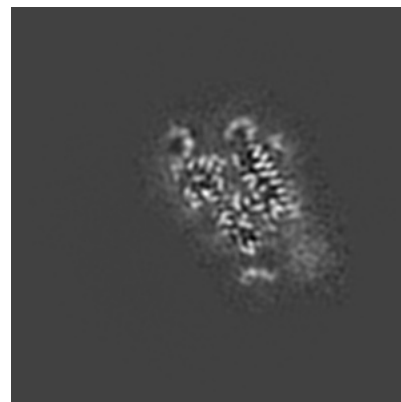
6.3.1 Primary map



X Index: 144

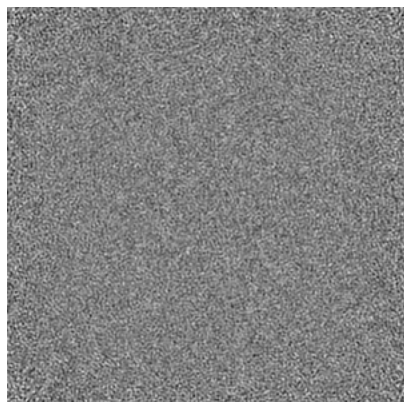


Y Index: 144

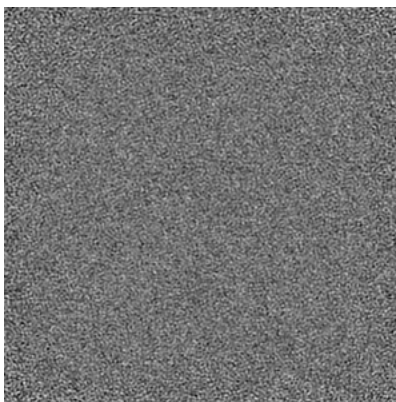


Z Index: 116

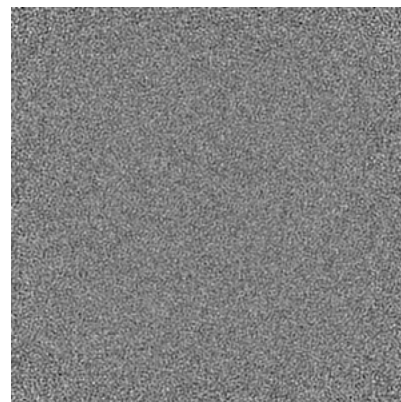
6.3.2 Raw map



X Index: 0



Y Index: 0

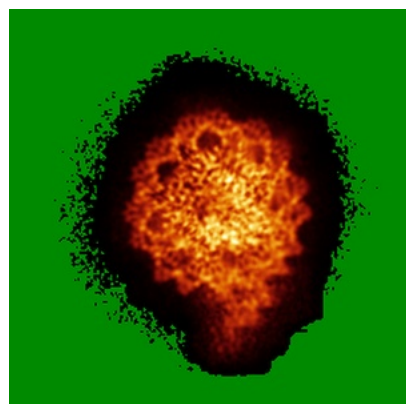


Z Index: 255

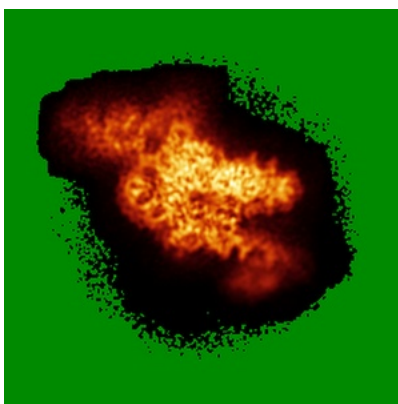
The images above show the largest variance slices of the map in three orthogonal directions.

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

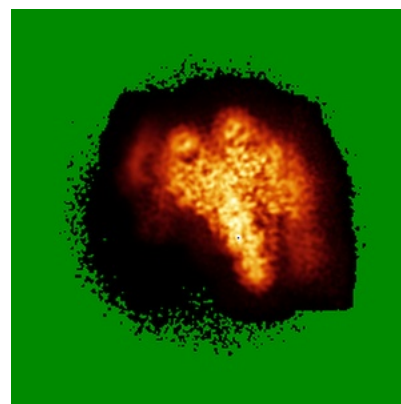
6.4.1 Primary map



X

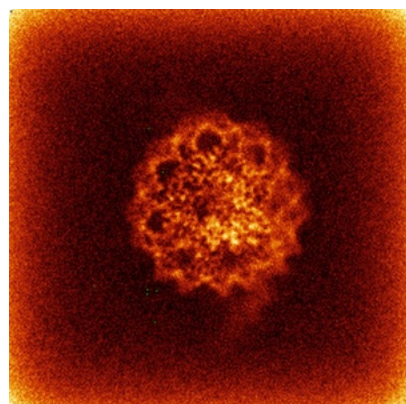


Y

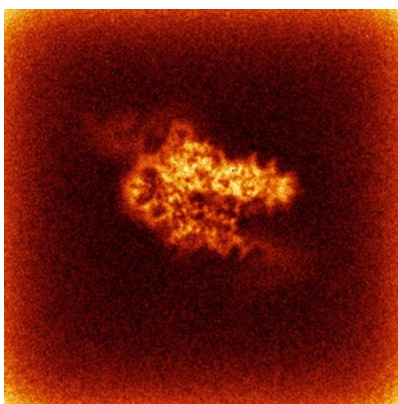


Z

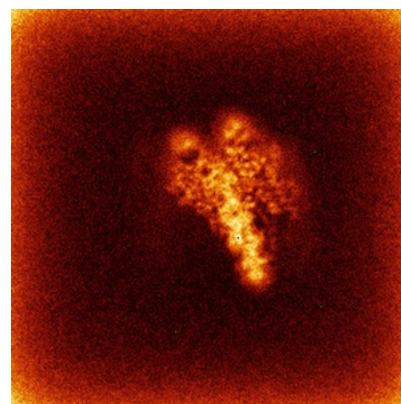
6.4.2 Raw map



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.

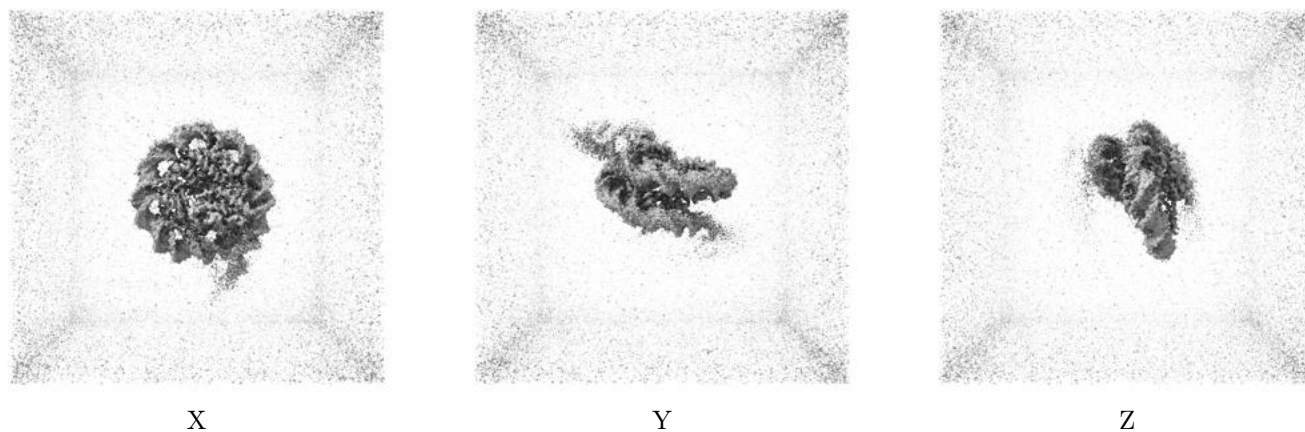
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.116. 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.

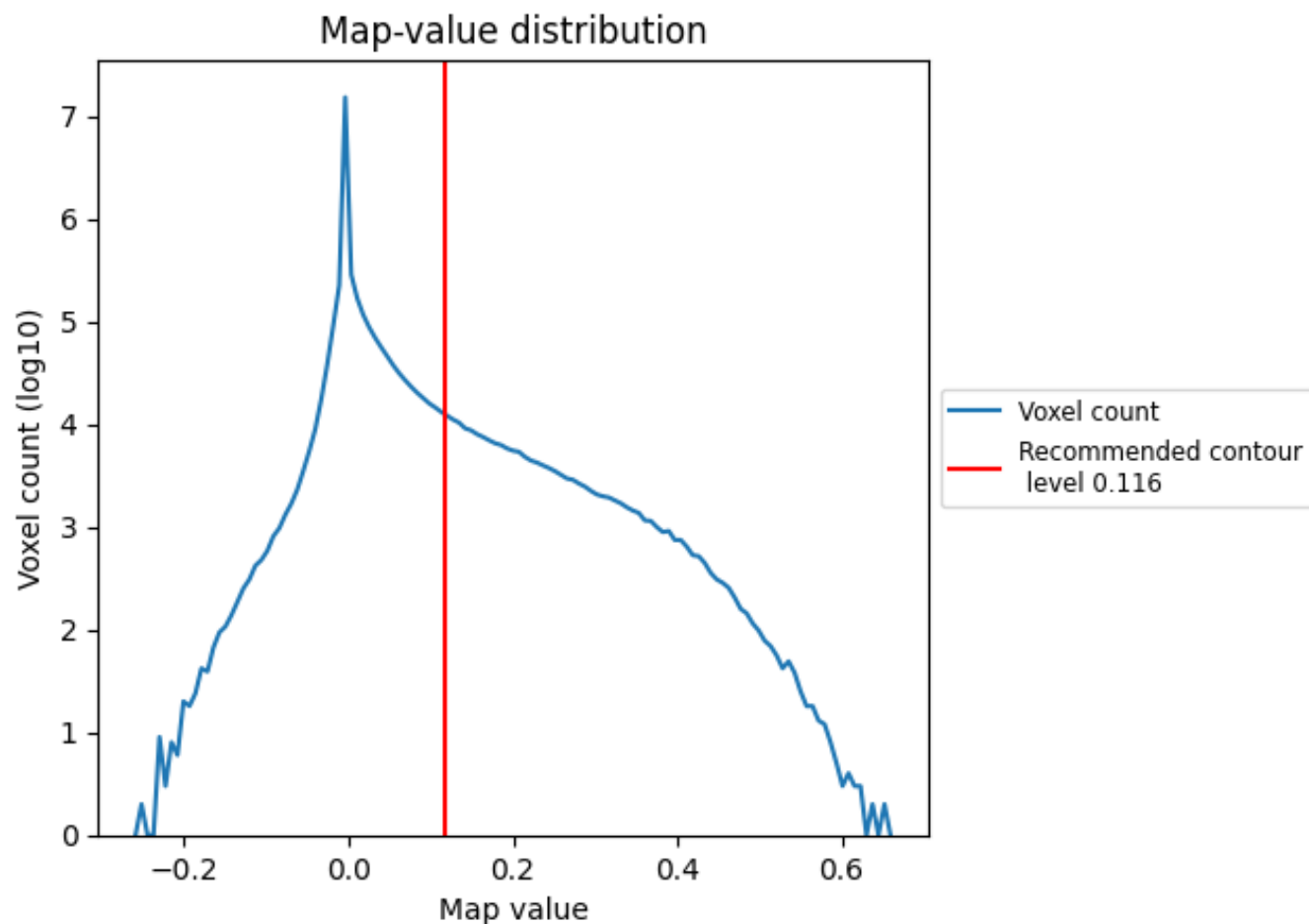
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

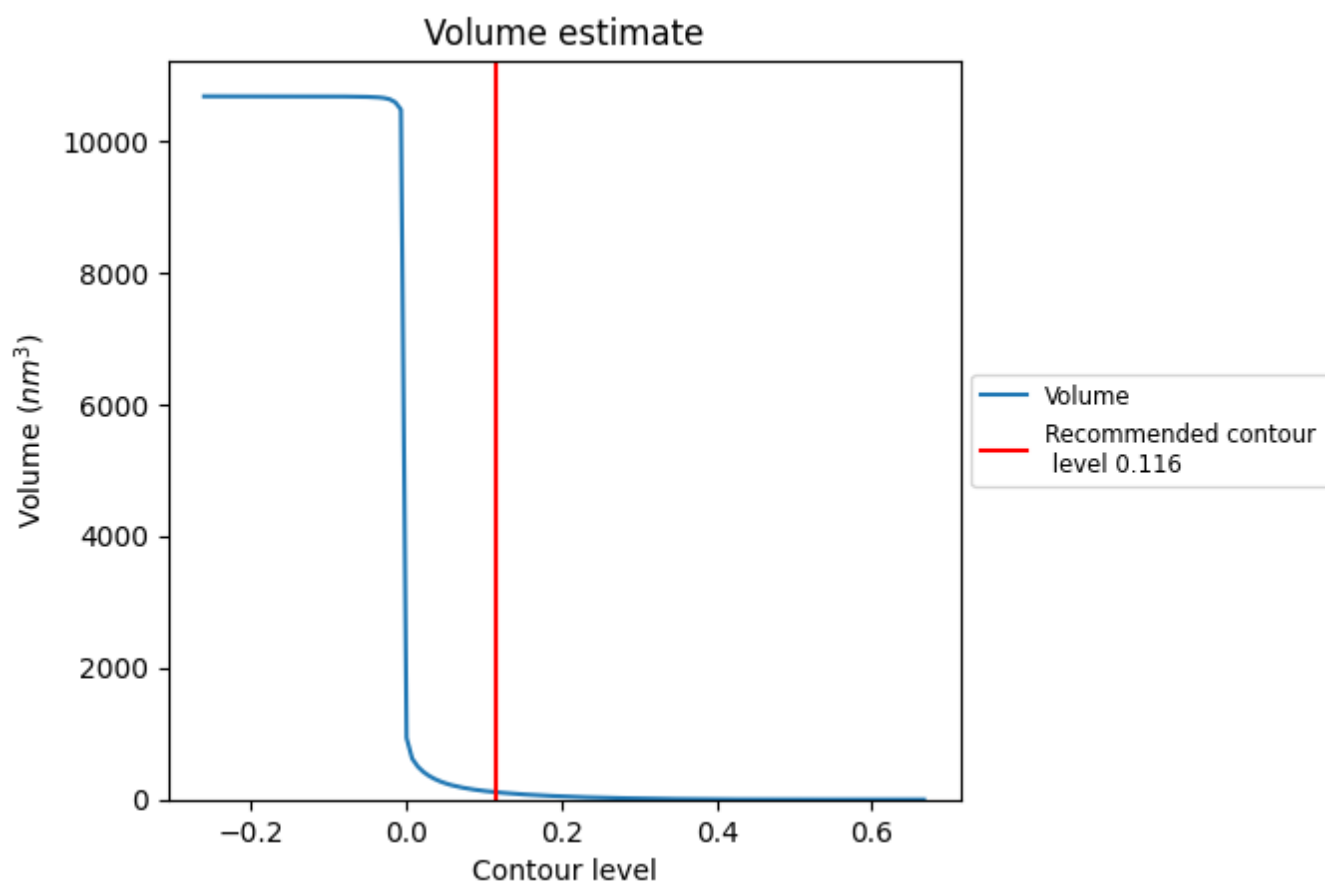
This section contains the results of statistical analysis of the map.

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.

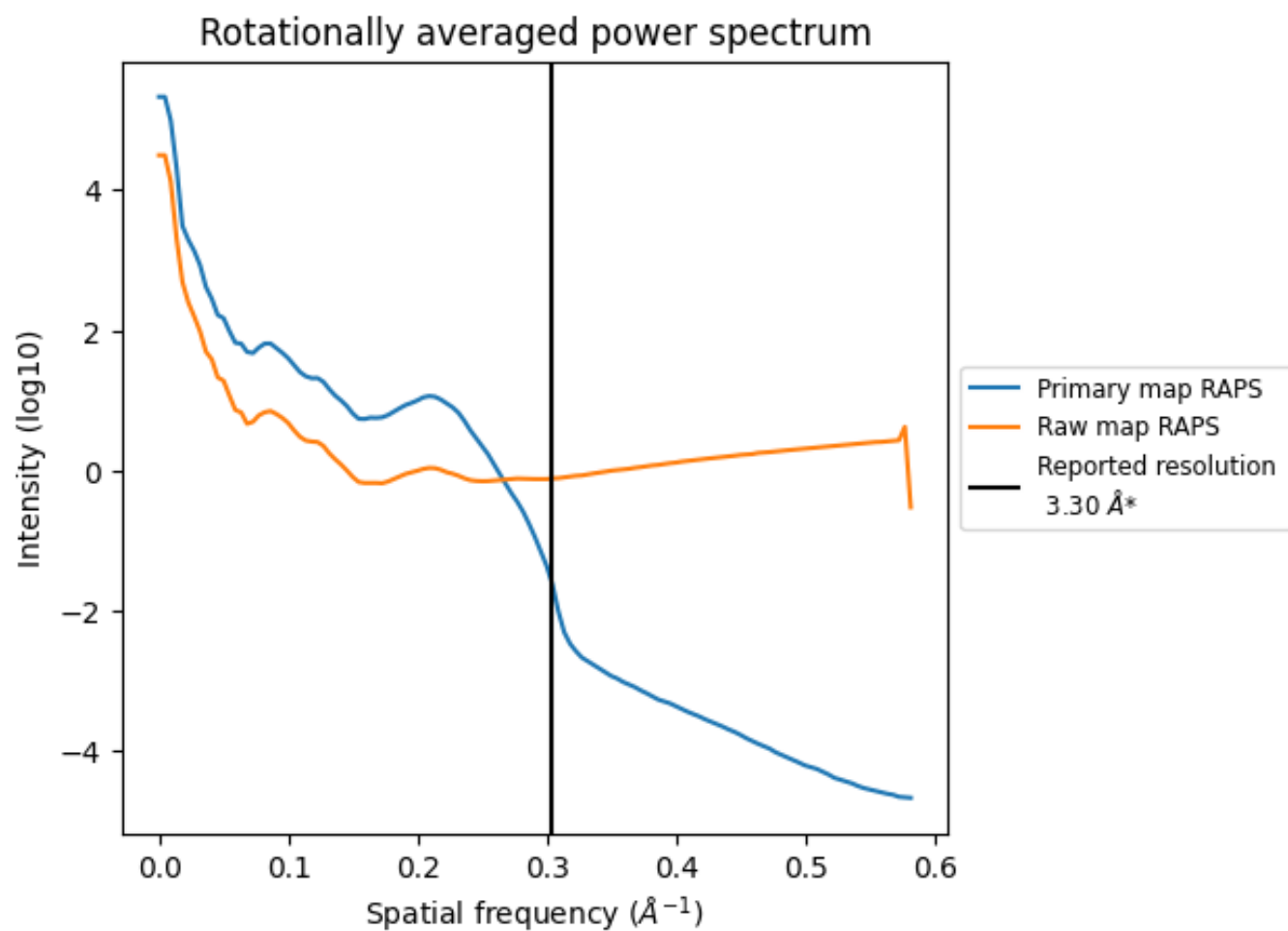
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 112 nm³; this corresponds to an approximate mass of 101 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 ⓘ

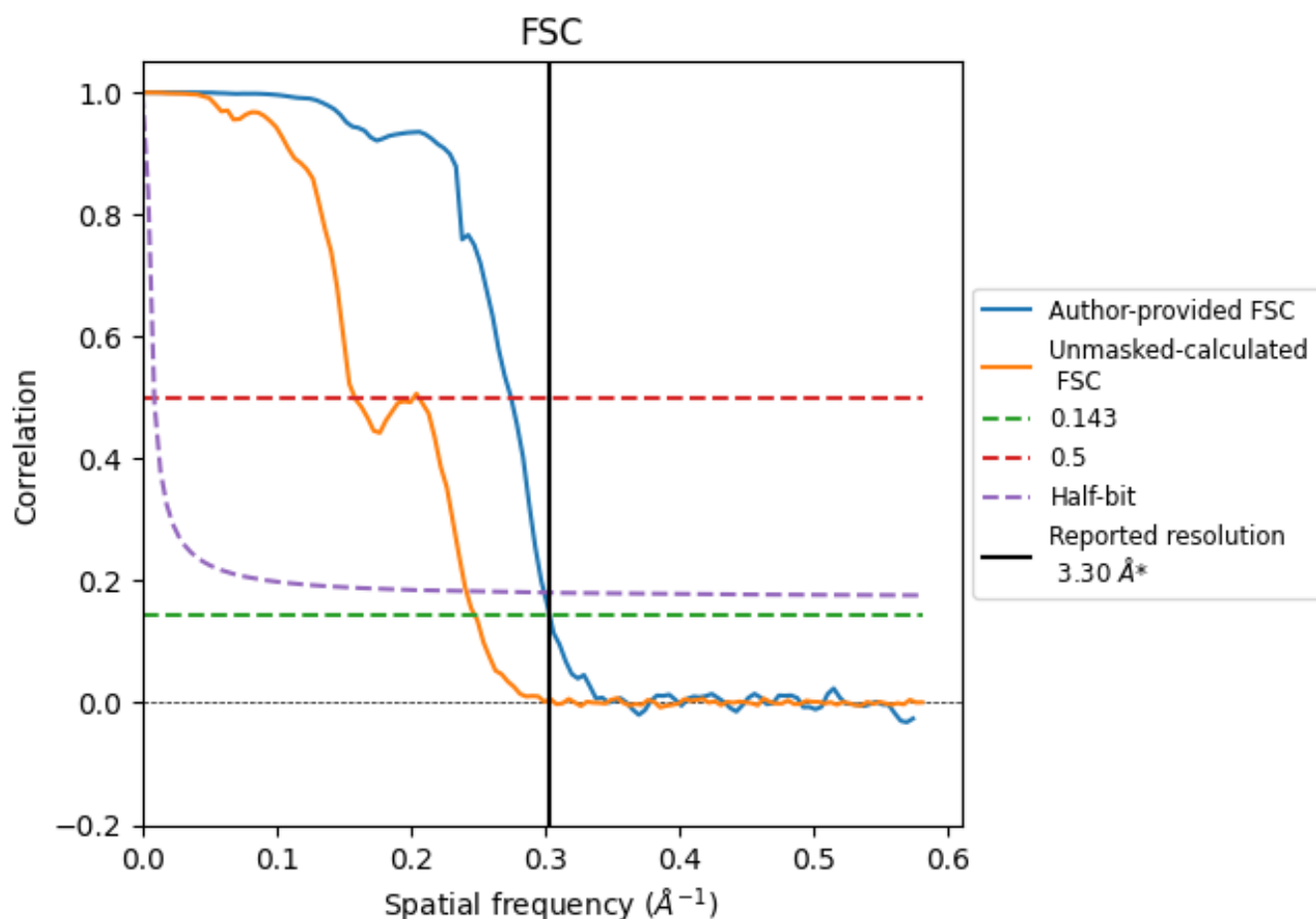


*Reported resolution corresponds to spatial frequency of 0.303 Å⁻¹

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 0.303 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

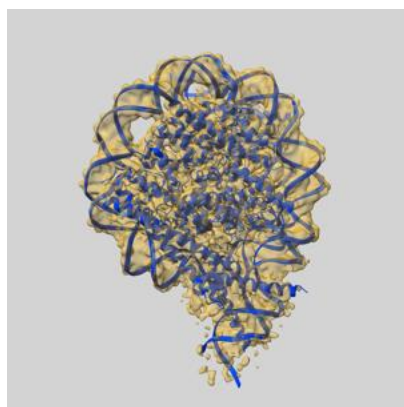
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.29	3.64	3.34
Unmasked-calculated*	4.03	6.31	4.14

*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 4.03 differs from the reported value 3.3 by more than 10 %

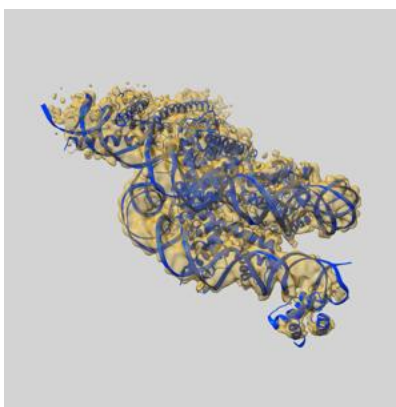
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-17183 and PDB model 8OTS. Per-residue inclusion information can be found in section [3](#) on page [9](#).

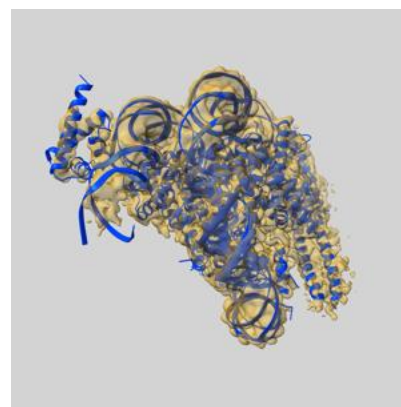
9.1 Map-model overlay [i](#)



X



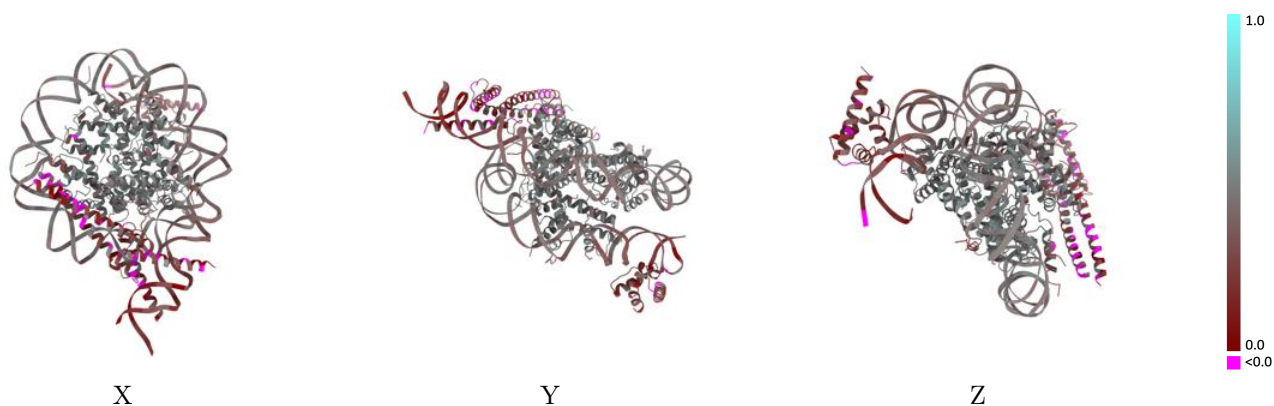
Y



Z

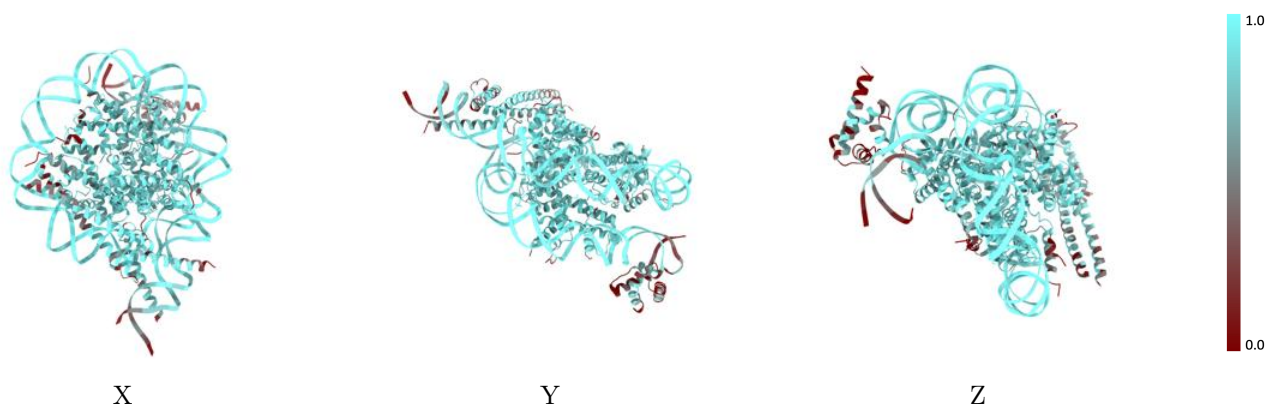
The images above show the 3D surface view of the map at the recommended contour level 0.116 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



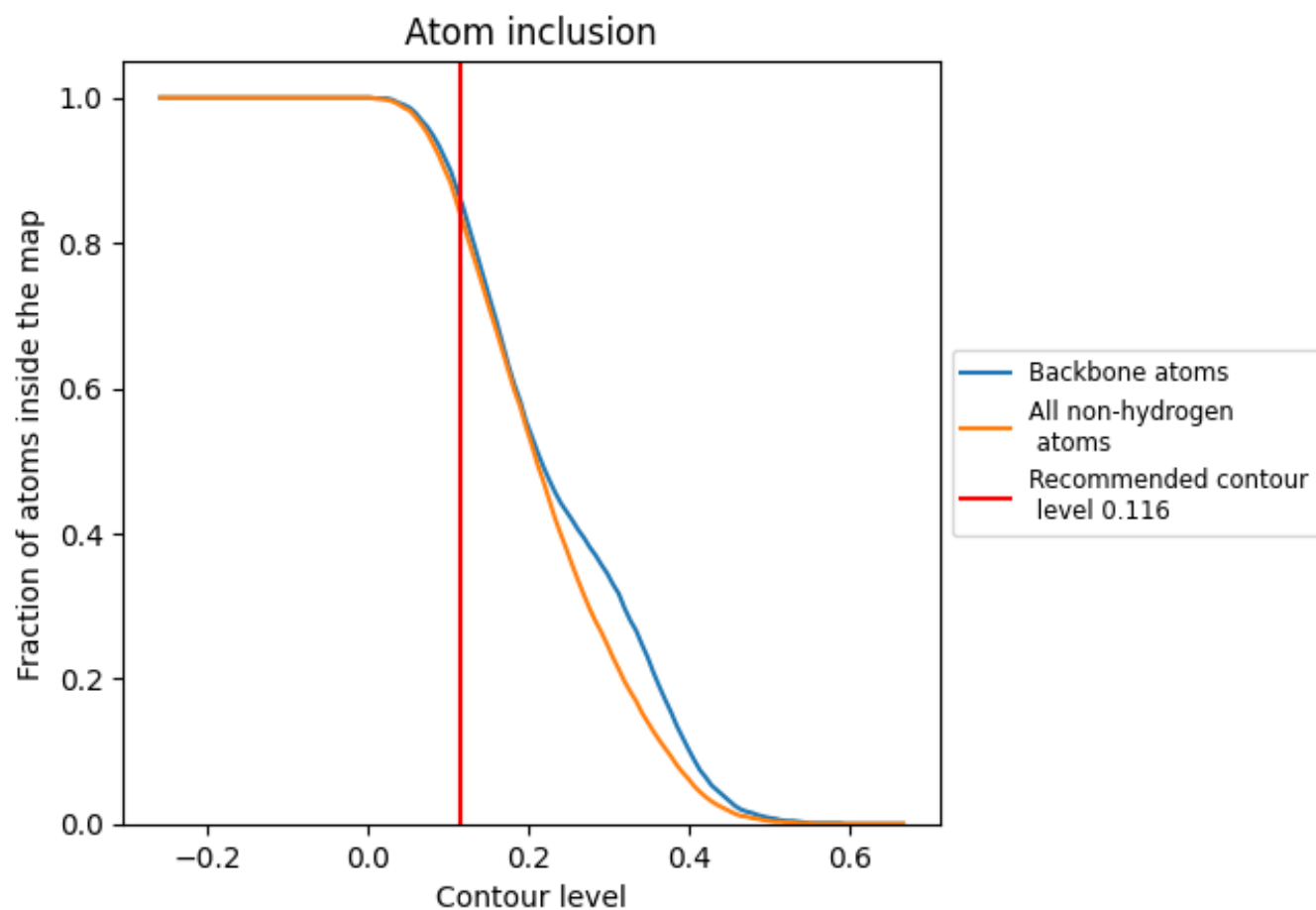
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.116).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% 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 (0.116) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8350	 0.4040
A	 0.8290	 0.4680
B	 0.8960	 0.4910
C	 0.8010	 0.4530
D	 0.8750	 0.4710
E	 0.8190	 0.4730
F	 0.8840	 0.4760
G	 0.8180	 0.4430
H	 0.8830	 0.4500
I	 0.8780	 0.3870
J	 0.8700	 0.3750
K	 0.3830	 0.2220
M	 0.6770	 0.1740
N	 0.7190	 0.1800

