



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 04:02 AM UTC

PDB ID : 9GXA / pdb_00009gxa
EMDB ID : EMD-51645
Title : CENP-A/H4 di-tetrasome assembled on alpha-satellite DNA.
Authors : Ali-Ahmad, A.; Sekulic, N.
Deposited on : 2024-09-29
Resolution : 4.01 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

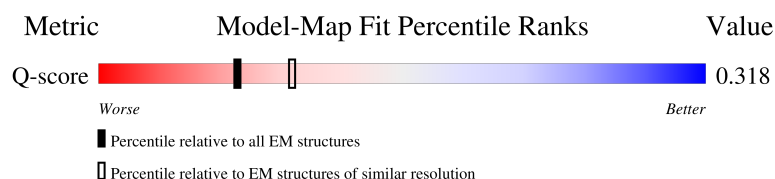
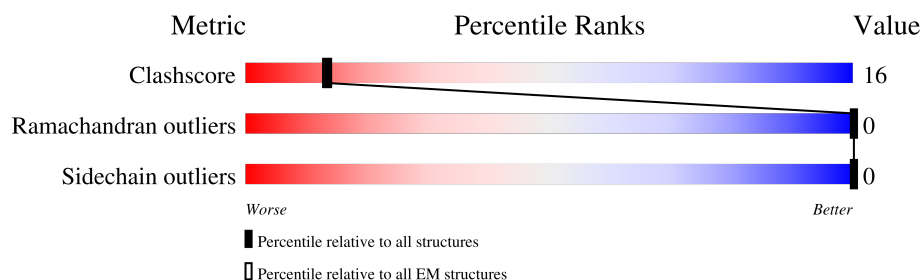
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.01 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	6765 (3.51 - 4.51)

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	17% 38% 13% 49%
1	C	139	45% 11% 45%
1	E	139	35% 21% 45%
1	G	139	10% 36% 19% 45%

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Mol	Chain	Length	Quality of chain
2	B	102	25%51%13%36%
2	D	102	50%18%32%
2	F	102	45%25%30%
2	H	102	11%42%25%33%
3	I	147	26%50%24%
4	J	147	22%54%24%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9215 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-like centromeric protein A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	71	Total	C	N	O	S	0	0
			575	375	104	95	1		
1	C	77	Total	C	N	O	S	0	0
			628	409	116	102	1		
1	E	77	Total	C	N	O	S	0	0
			628	409	116	102	1		
1	G	76	Total	C	N	O	S	0	0
			621	405	115	100	1		

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	65	Total	C	N	O	S	0	0
			516	328	98	89	1		
2	D	69	Total	C	N	O	S	0	0
			549	347	107	94	1		
2	F	71	Total	C	N	O	S	0	0
			565	355	110	99	1		
2	H	68	Total	C	N	O	S	0	0
			541	342	103	95	1		

- Molecule 3 is a DNA chain called 147 bp human alpha-satellite DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	112	Total	C	N	O	P	0	0
			2292	1097	412	671	112		

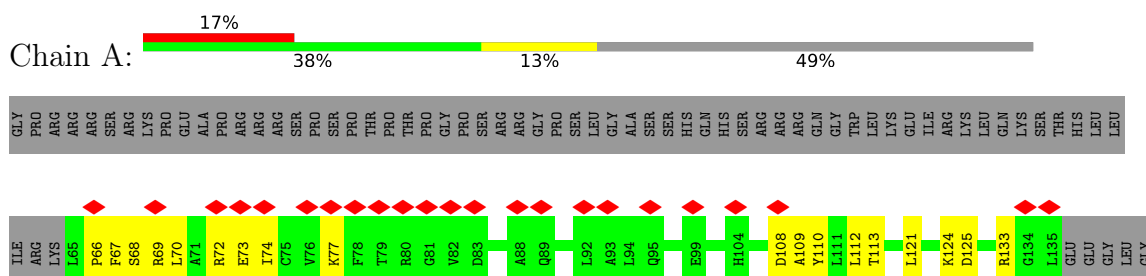
- Molecule 4 is a DNA chain called 147 bp alpha-satellite DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	J	112	Total	C	N	O	P	0	0
			2300	1100	415	673	112		

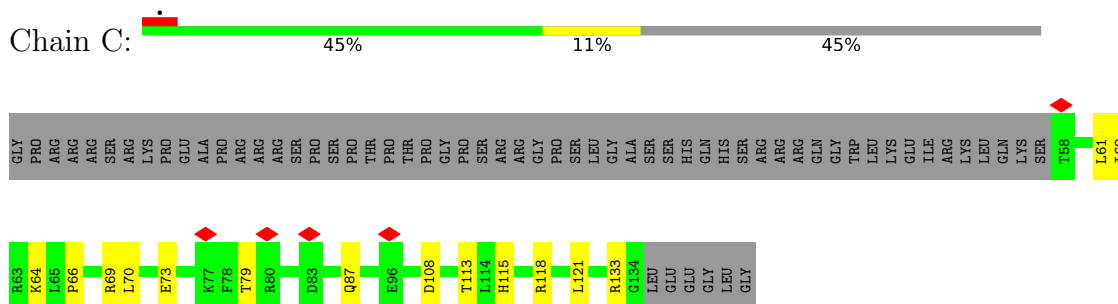
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

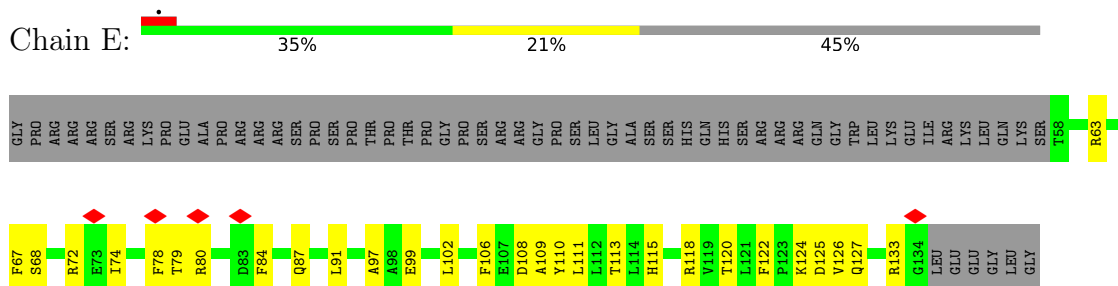
- Molecule 1: Histone H3-like centromeric protein A



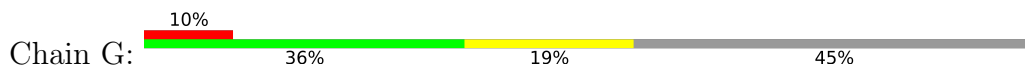
- Molecule 1: Histone H3-like centromeric protein A



- Molecule 1: Histone H3-like centromeric protein A



- Molecule 1: Histone H3-like centromeric protein A



Chain I:

A horizontal bar chart titled "Chain I:" showing the distribution of 60 categories. The bar is divided into three segments: green (26%), yellow (50%), and grey (24%). The categories are listed vertically below the bar, grouped by their segment color.

Category	Segment Color
C-7	Yellow
DC A-6	Grey
DC G-5	Yellow
DT T1	Yellow
DC G2	Grey
DC A3	Yellow
DA G4	Grey
DT T5	Yellow
DC G6	Grey
DC A7	Yellow
DC C13	Yellow
DC A14	Grey
DT T15	Yellow
DC C16	Yellow
DC A17	Grey
DT T18	Yellow
DC C21	Yellow
DC A22	Grey
DT A23	Yellow
DC A24	Grey
DT G25	Yellow
DC A26	Grey
DT T27	Yellow
DC G28	Grey
DT A29	Yellow
DC T30	Grey
DT T31	Yellow
DC G32	Green
DT T33	Yellow
DC G34	Green
DT A35	Yellow
DC G36	Green
DT A37	Yellow
DC T38	Yellow
DT A39	Green
DC T40	Yellow
DT T41	Green
DC G42	Green
DT A43	Yellow
DC T44	Yellow
DT T45	Yellow
DC A46	Yellow
DT T47	Yellow
DC A48	Yellow
DT T49	Yellow
DC G50	Yellow
DT C51	Yellow
DC T52	Yellow
DT T53	Yellow
DC T54	Yellow
DT T55	Yellow
DC A57	Green
DT T58	Yellow
DC A59	Yellow
DT T60	Yellow
DC G61	Yellow
DT A62	Yellow
DC DA	Grey
DT DC	Grey

Chain J:

DA
DT
DT
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DG
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DG
DA
DA
DA
DG
DT
T-62
C-61
A-60
A-54
A-53
G-52
G-51
G-44
A-43
A-42
G-41
C-40
A-39
T-38
T-37
C-36
T-35
C-34
A-33
T-27
T-26
C-25
T-24
T-23
T-22
G-21
T-20
G-19
A-18
T-17
G-16
A-15
T-14
G-13
G-12
A-11
T-8
T-7
C-4
T-3
C-2
A-1
C-1

G2
A3
G4
C5
T6
G7
A8
C9
C10
C11
T12
G13
C14
C15
T16
T17
T18
T19
G20
A21
T22
G23
G24
A25
G26
C27
A28
G29
C33
C34
A37
T38
A39
C40
A41
T44
T45
T46
G47
G48
T49
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	22500	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	57.5	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.339	Depositor
Minimum map value	-0.139	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.011	Depositor
Recommended contour level	0.108	Depositor
Map size ($\text{\AA}$)	295.2, 295.2, 295.2	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.82000005, 0.82000005, 0.82000005	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	0.15	0/587	0.31	0/794
1	C	0.46	0/641	0.56	0/866
1	E	0.33	0/641	0.48	0/866
1	G	0.15	0/634	0.36	0/856
2	B	0.15	0/521	0.31	0/700
2	D	0.19	0/554	0.34	0/742
2	F	0.20	0/570	0.34	0/764
2	H	0.15	0/546	0.34	0/733
3	I	0.21	0/2570	0.44	0/3963
4	J	0.20	0/2580	0.40	0/3981
All	All	0.23	0/9844	0.41	0/14265

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	575	0	591	20	0
1	C	628	0	653	24	0
1	E	628	0	653	30	0
1	G	621	0	646	22	0
2	B	516	0	558	11	0
2	D	549	0	595	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	565	0	605	28	0
2	H	541	0	581	23	0
3	I	2292	0	1267	62	0
4	J	2300	0	1268	79	0
All	All	9215	0	7417	266	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 16.

All (266) 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:112:LEU:HD11	1:A:125:ASP:OD1	1.32	1.29
1:A:112:LEU:CD1	1:A:125:ASP:OD1	1.93	1.16
1:A:113:THR:HG23	1:A:125:ASP:OD2	1.55	1.06
1:A:113:THR:CG2	1:A:125:ASP:OD2	2.12	0.96
1:C:121:LEU:HD11	2:D:43:VAL:CG2	1.99	0.92
1:C:121:LEU:HD11	2:D:43:VAL:HG22	1.58	0.84
1:C:113:THR:HG21	1:C:121:LEU:HD23	1.61	0.83
1:E:74:ILE:HD12	1:E:78:PHE:CD2	2.15	0.81
1:A:113:THR:CG2	1:A:125:ASP:CG	2.54	0.80
1:E:74:ILE:HD12	1:E:78:PHE:HD2	1.46	0.79
2:H:36:ARG:NH2	4:J:-44:DG:OP2	2.17	0.77
4:J:-17:DT:H4'	4:J:-16:DG:H5'	1.71	0.72
2:F:30:THR:HG21	3:I:17:DA:H5''	1.72	0.71
1:A:113:THR:HG22	1:A:125:ASP:CG	2.16	0.70
1:C:108:ASP:OD2	1:C:133:ARG:NH1	2.25	0.69
1:E:99:GLU:OE2	2:F:40:ARG:NH1	2.27	0.68
3:I:-36:DT:H2''	3:I:-35:DT:C5	2.30	0.67
1:C:121:LEU:HD11	2:D:43:VAL:HG21	1.77	0.67
1:A:112:LEU:HD12	1:A:125:ASP:OD1	1.91	0.67
1:G:61:LEU:HD23	2:H:36:ARG:HG3	1.77	0.67
1:E:79:THR:HG22	1:E:79:THR:O	1.96	0.66
1:E:80:ARG:O	1:E:80:ARG:HG3	1.95	0.65
2:F:30:THR:OG1	3:I:18:DT:OP1	2.13	0.65
3:I:13:DC:H2''	3:I:14:DA:C8	2.32	0.65
4:J:-23:DT:H2'	4:J:-22:DT:H71	1.80	0.64
1:C:121:LEU:CD1	2:D:43:VAL:HG21	2.29	0.62
2:D:72:TYR:HB3	2:F:75:HIS:CE1	2.35	0.61
4:J:-25:DC:H2''	4:J:-24:DT:H72	1.81	0.61
1:C:61:LEU:HD11	2:D:37:LEU:HA	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:43:VAL:HG21	2:B:46:ILE:HD11	1.82	0.60
2:H:26:ILE:HG23	2:H:55:ARG:HH21	1.67	0.60
3:I:-28:DT:H2''	3:I:-27:DG:C8	2.36	0.60
1:E:68:SER:HG	1:E:72:ARG:HH12	1.50	0.60
1:C:121:LEU:HD12	2:D:43:VAL:HG11	1.84	0.60
1:A:113:THR:HG23	1:A:125:ASP:CG	2.22	0.59
1:E:74:ILE:HD13	2:F:59:LYS:HE3	1.85	0.59
1:E:110:TYR:HB2	2:F:43:VAL:HG12	1.85	0.58
1:G:63:ARG:HE	3:I:48:DT:H4'	1.67	0.58
4:J:-26:DT:H2''	4:J:-25:DC:C6	2.39	0.58
3:I:46:DG:H2''	3:I:47:DT:H71	1.84	0.58
2:H:32:PRO:HB2	2:H:36:ARG:HH21	1.69	0.57
4:J:-24:DT:H2'	4:J:-23:DT:H71	1.87	0.57
1:A:124:LYS:HG3	1:C:115:HIS:NE2	2.20	0.56
3:I:30:DT:H2'	3:I:31:DT:H71	1.87	0.56
1:G:120:THR:HG22	2:H:45:ARG:HD3	1.87	0.56
1:A:74:ILE:HD12	1:A:77:LYS:HD2	1.86	0.56
2:H:57:VAL:HA	2:H:60:VAL:HG12	1.88	0.56
4:J:15:DC:H2''	4:J:16:DT:C5	2.40	0.56
3:I:-7:DC:H2''	3:I:-6:DA:C8	2.41	0.55
3:I:-41:DT:H2''	3:I:-40:DG:H5''	1.88	0.55
3:I:-5:DG:N2	4:J:6:DT:O2	2.39	0.55
1:G:123:PRO:HG2	2:H:49:LEU:HD23	1.88	0.55
2:H:24:ASP:HA	2:H:27:GLN:HG2	1.88	0.55
3:I:34:DG:H2''	3:I:35:DA:C8	2.41	0.55
2:H:52:GLU:OE1	2:H:52:GLU:N	2.34	0.55
3:I:2:DG:H2''	3:I:3:DA:C8	2.42	0.55
3:I:33:DT:H2''	3:I:34:DG:C8	2.42	0.55
1:G:66:PRO:HA	1:G:69:ARG:HG2	1.89	0.54
1:E:84:PHE:HE2	2:F:70:VAL:HG13	1.72	0.54
3:I:36:DG:H2''	3:I:37:DA:C8	2.41	0.54
4:J:37:DA:H2''	4:J:38:DT:H71	1.88	0.54
1:G:74:ILE:HG23	2:H:66:ILE:HD13	1.89	0.54
2:H:25:ASN:OD1	2:H:26:ILE:N	2.41	0.54
2:D:46:ILE:HD11	2:D:50:ILE:HG21	1.88	0.54
1:E:126:VAL:HG21	2:F:50:ILE:HD11	1.89	0.54
3:I:58:DT:H2''	3:I:59:DA:N7	2.22	0.54
4:J:46:DT:H2''	4:J:47:DG:C8	2.42	0.54
4:J:18:DT:H2''	4:J:19:DT:H71	1.89	0.54
2:H:39:ARG:O	2:H:39:ARG:NH1	2.36	0.54
1:C:121:LEU:HB2	2:D:46:ILE:HD12	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:102:LEU:HB3	2:H:37:LEU:HD21	1.91	0.53
1:C:121:LEU:CD1	2:D:43:VAL:CG2	2.79	0.53
3:I:52:DC:H2'	3:I:53:DT:H71	1.90	0.53
1:G:72:ARG:HG2	1:G:86:TRP:HE1	1.73	0.53
4:J:38:DT:H2''	4:J:39:DA:C8	2.44	0.53
1:G:74:ILE:HD12	1:G:77:LYS:HG2	1.90	0.53
4:J:16:DT:H2''	4:J:17:DT:H5''	1.90	0.53
3:I:24:DA:H2''	3:I:25:DG:C8	2.43	0.52
1:C:61:LEU:HD12	2:D:36:ARG:HG2	1.90	0.52
1:E:97:ALA:HB2	2:F:90:LEU:HD21	1.91	0.52
3:I:51:DC:H2''	3:I:52:DC:H5''	1.91	0.52
4:J:24:DG:H2''	4:J:25:DA:C8	2.45	0.52
1:G:78:PHE:HB3	1:G:80:ARG:NE	2.24	0.52
1:E:87:GLN:HG3	3:I:7:DA:OP1	2.09	0.51
3:I:28:DT:H2''	3:I:29:DA:C8	2.46	0.51
3:I:-13:DC:H2''	3:I:-12:DA:C8	2.45	0.51
1:G:119:VAL:N	4:J:-34:DC:OP1	2.43	0.51
3:I:55:DT:H2''	3:I:56:DT:H72	1.92	0.51
1:C:73:GLU:HG2	2:D:26:ILE:HD12	1.92	0.51
3:I:14:DA:H2''	3:I:15:DT:H72	1.92	0.51
2:B:39:ARG:NH1	2:B:44:LYS:O	2.44	0.51
4:J:25:DA:H2''	4:J:26:DG:C8	2.46	0.51
1:A:67:PHE:CD1	2:B:28:GLY:HA3	2.45	0.50
1:C:87:GLN:HG3	4:J:7:DG:OP1	2.10	0.50
1:C:62:ILE:HD11	2:D:29:ILE:HG12	1.92	0.50
4:J:-40:DC:H2''	4:J:-39:DA:C8	2.47	0.50
2:H:47:SER:HA	3:I:38:DA:O5'	2.11	0.50
2:F:51:TYR:HB3	2:F:55:ARG:HH12	1.77	0.50
1:E:63:ARG:HH12	3:I:17:DA:H5'	1.76	0.50
4:J:-42:DA:C8	4:J:-42:DA:H5'	2.47	0.50
1:C:69:ARG:HD3	2:D:26:ILE:HG12	1.93	0.49
4:J:-2:DC:H2''	4:J:-1:DA:C8	2.47	0.49
4:J:13:DG:H2''	4:J:14:DC:C6	2.47	0.49
1:A:113:THR:HG21	1:A:121:LEU:HD13	1.94	0.49
1:E:68:SER:HB2	1:E:91:LEU:HD21	1.93	0.49
3:I:-23:DC:H1'	3:I:-22:DA:H5'	1.94	0.49
1:G:62:ILE:HD11	1:G:95:GLN:OE1	2.12	0.49
1:G:62:ILE:HG22	2:H:33:ALA:HB1	1.93	0.49
2:D:70:VAL:HA	2:D:73:THR:HG22	1.94	0.49
1:E:68:SER:OG	1:E:72:ARG:NH1	2.34	0.49
3:I:5:DT:H2''	3:I:6:DG:C8	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:ALA:O	1:A:113:THR:HG23	2.13	0.49
2:B:48:GLY:N	4:J:38:DT:OP1	2.34	0.49
2:D:92:ARG:HG2	2:F:92:ARG:HH12	1.76	0.49
3:I:21:DC:H2'	3:I:22:DA:C8	2.47	0.49
2:B:62:LEU:HA	2:B:65:VAL:HG22	1.94	0.48
3:I:-26:DC:H2''	3:I:-25:DT:C5	2.48	0.48
4:J:-25:DC:H2''	4:J:-24:DT:C7	2.43	0.48
4:J:-14:DT:H2''	4:J:-13:DG:C8	2.48	0.48
4:J:3:DA:H2''	4:J:4:DG:C8	2.47	0.48
4:J:17:DT:H2'	4:J:18:DT:H71	1.95	0.48
1:C:79:THR:HG21	2:D:70:VAL:HG21	1.95	0.48
4:J:-37:DT:H2''	4:J:-36:DC:C6	2.48	0.48
1:E:74:ILE:CD1	1:E:78:PHE:CD2	2.93	0.48
1:C:70:LEU:O	1:C:73:GLU:HG3	2.14	0.47
2:D:83:ALA:HA	2:D:86:VAL:HG22	1.97	0.47
3:I:5:DT:H2''	3:I:6:DG:N7	2.29	0.47
3:I:-45:DA:H2''	3:I:-44:DA:C8	2.50	0.47
3:I:23:DA:H2''	3:I:24:DA:C8	2.49	0.47
3:I:15:DT:H2''	3:I:16:DC:C5	2.49	0.47
4:J:-12:DG:H2''	4:J:-11:DA:H8	1.79	0.47
3:I:-39:DT:H2''	3:I:-38:DA:C8	2.50	0.46
4:J:33:DC:H2''	4:J:34:DC:C5	2.50	0.46
1:E:79:THR:O	1:E:79:THR:CG2	2.63	0.46
3:I:-36:DT:H2''	3:I:-35:DT:C7	2.45	0.46
3:I:46:DG:H2''	3:I:47:DT:C7	2.44	0.46
4:J:-34:DC:H2''	4:J:-33:DA:C8	2.50	0.46
4:J:27:DC:H2''	4:J:28:DA:C8	2.50	0.46
4:J:28:DA:H2''	4:J:29:DG:H8	1.80	0.46
4:J:-35:DT:H6	4:J:-35:DT:H2'	1.63	0.46
1:A:108:ASP:OD2	1:A:133:ARG:NH1	2.49	0.46
4:J:5:DC:H2''	4:J:6:DT:C5	2.51	0.46
4:J:2:DG:H2''	4:J:3:DA:C8	2.51	0.46
1:A:66:PRO:O	1:A:69:ARG:HG2	2.15	0.46
2:D:75:HIS:HA	2:F:88:TYR:CE2	2.50	0.46
2:F:31:LYS:HG3	2:F:55:ARG:HH22	1.80	0.46
3:I:1:DT:H2''	3:I:2:DG:C8	2.51	0.46
3:I:27:DA:H2''	3:I:28:DT:H71	1.98	0.46
4:J:22:DT:C4	4:J:23:DG:C6	3.04	0.46
2:F:47:SER:HA	4:J:-24:DT:H5''	1.98	0.46
2:H:26:ILE:HG23	2:H:55:ARG:NH2	2.30	0.46
4:J:0:DC:H2''	4:J:1:DA:C8	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:118:ARG:HH21	1:E:122:PHE:HB2	1.79	0.45
3:I:45:DC:H2''	3:I:46:DG:N7	2.31	0.45
1:E:106:PHE:HD2	2:F:38:ALA:HA	1.80	0.45
4:J:-26:DT:H2''	4:J:-25:DC:C5	2.52	0.45
4:J:-52:DG:C2	4:J:-51:DG:C5	3.04	0.45
4:J:-13:DG:H2''	4:J:-12:DG:C8	2.52	0.45
2:F:72:TYR:HE1	2:F:88:TYR:HB3	1.82	0.45
3:I:3:DA:H2''	3:I:4:DG:C8	2.52	0.45
3:I:49:DT:H2''	3:I:50:DG:C8	2.52	0.45
4:J:-8:DT:H2''	4:J:-7:DT:C6	2.52	0.45
2:F:51:TYR:O	2:F:55:ARG:HG3	2.17	0.45
1:G:67:PHE:O	1:G:70:LEU:HG	2.17	0.45
1:G:110:TYR:HA	1:G:113:THR:HG22	1.97	0.45
1:E:102:LEU:HD11	2:F:58:LEU:HD13	1.99	0.45
4:J:-39:DA:H1'	4:J:-38:DT:H5'	1.98	0.45
4:J:-16:DG:H1'	4:J:-15:DA:C8	2.52	0.45
4:J:-13:DG:H2''	4:J:-12:DG:H8	1.82	0.45
4:J:38:DT:H2''	4:J:39:DA:H8	1.81	0.45
1:C:64:LYS:HE2	1:C:64:LYS:HB3	1.69	0.44
3:I:-18:DA:H2''	3:I:-17:DA:C8	2.52	0.44
1:G:70:LEU:HA	1:G:73:GLU:OE2	2.17	0.44
1:E:80:ARG:O	1:E:80:ARG:CG	2.65	0.44
1:E:115:HIS:CE1	1:G:124:LYS:HE3	2.53	0.44
3:I:-9:DT:H2'	3:I:-8:DT:H71	1.99	0.44
2:H:48:GLY:N	3:I:38:DA:OP2	2.35	0.44
2:D:75:HIS:CE1	2:F:76:ALA:HB2	2.53	0.44
4:J:19:DT:H2''	4:J:20:DG:H8	1.82	0.44
2:F:80:THR:HG21	4:J:-3:DT:H5''	2.00	0.44
1:A:68:SER:O	1:A:72:ARG:HG2	2.17	0.44
1:E:109:ALA:O	1:E:113:THR:HG23	2.18	0.44
4:J:-54:DA:H2''	4:J:-53:DA:C8	2.53	0.44
4:J:-20:DT:N3	4:J:-19:DG:O6	2.50	0.43
4:J:21:DA:H2'	4:J:22:DT:H71	2.00	0.43
1:A:73:GLU:O	1:A:77:LYS:NZ	2.50	0.43
2:F:30:THR:HG23	2:F:33:ALA:H	1.83	0.43
3:I:44:DC:H2''	3:I:45:DC:C6	2.53	0.43
4:J:12:DT:H2''	4:J:13:DG:C8	2.53	0.43
4:J:-8:DT:H2''	4:J:-7:DT:C5	2.54	0.43
1:E:79:THR:HG21	2:F:70:VAL:HG11	2.00	0.43
3:I:37:DA:H2''	3:I:38:DA:C8	2.53	0.43
4:J:-61:DC:H2''	4:J:-60:DA:C8	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:GLU:O	1:A:77:LYS:HG3	2.18	0.43
3:I:-32:DA:H2''	3:I:-31:DA:H8	1.83	0.43
4:J:20:DG:H4'	4:J:21:DA:OP1	2.17	0.43
2:D:73:THR:HG21	2:D:81:VAL:HG22	2.01	0.43
3:I:-44:DA:H2''	3:I:-43:DA:C8	2.54	0.43
4:J:-22:DT:C2	4:J:-21:DG:N7	2.87	0.43
2:F:83:ALA:HA	2:F:86:VAL:HG22	2.00	0.43
1:G:100:ALA:HA	1:G:103:VAL:HG12	2.01	0.43
4:J:9:DA:C4	4:J:10:DC:C5	3.07	0.43
4:J:39:DA:C4	4:J:40:DC:C5	3.07	0.43
4:J:40:DC:C2	4:J:41:DA:N7	2.87	0.42
4:J:44:DT:H4'	4:J:45:DT:OP2	2.16	0.42
1:C:121:LEU:HD12	2:D:46:ILE:CD1	2.49	0.42
2:D:75:HIS:HE1	2:F:76:ALA:HB2	1.83	0.42
1:E:67:PHE:CE1	2:F:62:LEU:HD11	2.54	0.42
3:I:-41:DT:H2''	3:I:-40:DG:H8	1.83	0.42
3:I:-26:DC:H2''	3:I:-25:DT:C7	2.49	0.42
3:I:-6:DA:H2''	3:I:-5:DG:C8	2.54	0.42
4:J:-12:DG:C2	4:J:-11:DA:C5	3.07	0.42
1:C:66:PRO:HA	1:C:69:ARG:HG2	2.01	0.42
3:I:26:DA:C8	3:I:26:DA:H5'	2.55	0.42
1:A:110:TYR:CE1	2:B:43:VAL:HA	2.54	0.42
2:H:46:ILE:HG23	2:H:50:ILE:HG21	2.01	0.42
4:J:0:DC:H2''	4:J:1:DA:H8	1.83	0.42
2:B:48:GLY:H	4:J:38:DT:P	2.40	0.42
2:D:63:GLU:O	2:D:67:ARG:HB2	2.19	0.42
4:J:-4:DC:H2'	4:J:-3:DT:H72	2.01	0.42
4:J:33:DC:H2''	4:J:34:DC:C6	2.54	0.42
2:H:60:VAL:HA	2:H:63:GLU:HG2	2.01	0.42
4:J:-44:DG:H2''	4:J:-43:DA:H8	1.84	0.42
2:B:54:THR:HA	2:B:57:VAL:HG12	2.00	0.42
2:F:75:HIS:CD2	2:F:75:HIS:C	2.98	0.42
3:I:23:DA:H2''	3:I:24:DA:H8	1.85	0.42
4:J:24:DG:H4'	4:J:25:DA:H5'	2.02	0.42
1:E:111:LEU:HD23	1:E:111:LEU:HA	1.85	0.42
2:F:79:LYS:HE2	2:F:79:LYS:HB2	1.96	0.42
1:G:71:ALA:HA	1:G:74:ILE:HG22	2.02	0.42
1:C:113:THR:HB	1:C:118:ARG:O	2.20	0.41
3:I:-45:DA:H2''	3:I:-44:DA:N7	2.35	0.41
2:B:57:VAL:HA	2:B:60:VAL:HG22	2.02	0.41
2:F:36:ARG:NH2	3:I:17:DA:OP1	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:-48:DC:H2''	3:I:-47:DC:C5	2.54	0.41
2:H:58:LEU:O	2:H:62:LEU:HD23	2.20	0.41
3:I:-47:DC:H2''	3:I:-46:DA:N7	2.36	0.41
3:I:47:DT:H2''	3:I:48:DT:H72	2.01	0.41
1:G:85:ASN:O	2:H:81:VAL:HG22	2.20	0.41
4:J:-61:DC:H2''	4:J:-60:DA:H8	1.86	0.41
1:E:108:ASP:OD2	1:E:133:ARG:NH2	2.47	0.41
1:E:124:LYS:HA	1:E:127:GLN:OE1	2.20	0.41
1:C:121:LEU:CD1	2:D:43:VAL:CG1	2.99	0.41
2:D:46:ILE:HD11	2:D:50:ILE:HD13	2.03	0.41
1:G:78:PHE:CZ	2:H:66:ILE:HG23	2.55	0.41
4:J:20:DG:C6	4:J:21:DA:C6	3.08	0.41
1:C:113:THR:CG2	1:C:121:LEU:HD23	2.42	0.41
2:D:72:TYR:CE2	2:D:88:TYR:HB3	2.55	0.41
4:J:-22:DT:C2	4:J:-21:DG:C8	3.09	0.41
4:J:-42:DA:C6	4:J:-41:DG:C6	3.08	0.41
4:J:5:DC:H2''	4:J:6:DT:C6	2.56	0.41
1:E:120:THR:HG22	3:I:28:DT:OP1	2.21	0.40
3:I:27:DA:C2'	3:I:28:DT:H71	2.51	0.40
4:J:-27:DT:H2''	4:J:-26:DT:C6	2.56	0.40
1:G:119:VAL:HG22	4:J:-34:DC:P	2.62	0.40
2:H:30:THR:OG1	2:H:32:PRO:HD2	2.20	0.40
3:I:-31:DA:H2''	3:I:-30:DA:H8	1.86	0.40
4:J:10:DC:C2	4:J:11:DA:N7	2.89	0.40
2:B:38:ALA:HB1	2:B:43:VAL:HG21	2.04	0.40
4:J:-22:DT:H2''	4:J:-21:DG:H8	1.86	0.40
4:J:20:DG:H2''	4:J:21:DA:H8	1.85	0.40
2:B:59:LYS:O	2:B:62:LEU:HG	2.21	0.40
1:E:113:THR:HG22	1:E:125:ASP:OD2	2.21	0.40
4:J:23:DG:H2''	4:J:24:DG:C8	2.56	0.40
1:A:70:LEU:O	1:A:73:GLU:HG2	2.21	0.40
4:J:-21:DG:C5	4:J:-20:DT:C4	3.09	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	69/139 (50%)	69 (100%)	0	0	100	100
1	C	75/139 (54%)	73 (97%)	2 (3%)	0	100	100
1	E	75/139 (54%)	74 (99%)	1 (1%)	0	100	100
1	G	74/139 (53%)	73 (99%)	1 (1%)	0	100	100
2	B	63/102 (62%)	63 (100%)	0	0	100	100
2	D	67/102 (66%)	65 (97%)	2 (3%)	0	100	100
2	F	69/102 (68%)	67 (97%)	2 (3%)	0	100	100
2	H	66/102 (65%)	65 (98%)	1 (2%)	0	100	100
All	All	558/964 (58%)	549 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	58/117 (50%)	58 (100%)	0	100	100
1	C	64/117 (55%)	64 (100%)	0	100	100
1	E	64/117 (55%)	64 (100%)	0	100	100
1	G	63/117 (54%)	63 (100%)	0	100	100
2	B	54/78 (69%)	54 (100%)	0	100	100
2	D	57/78 (73%)	57 (100%)	0	100	100
2	F	59/78 (76%)	59 (100%)	0	100	100
2	H	57/78 (73%)	57 (100%)	0	100	100
All	All	476/780 (61%)	476 (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 (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	85	ASN
1	E	115	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](#)

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.

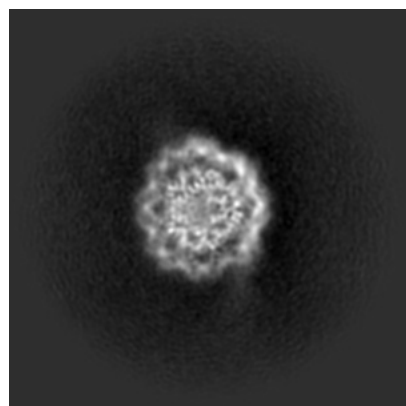
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-51645. These allow visual inspection of the internal detail of the map and identification of artifacts.

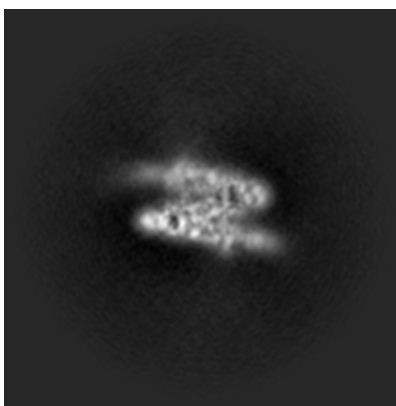
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



X

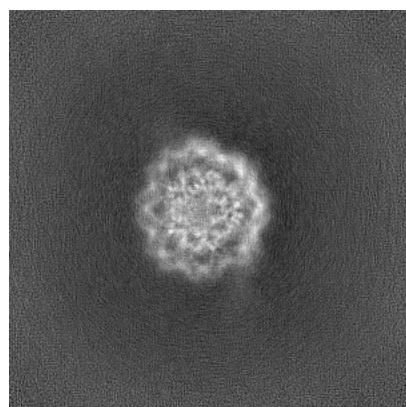


Y

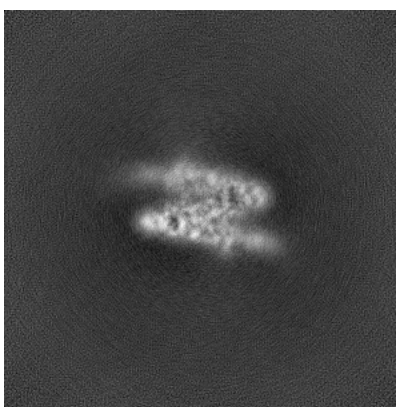


Z

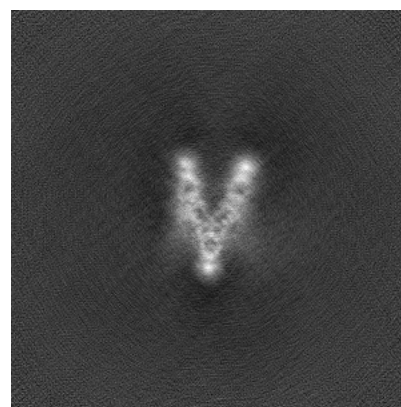
6.1.2 Raw map



X



Y

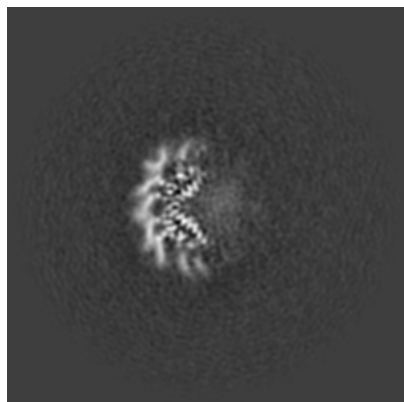


Z

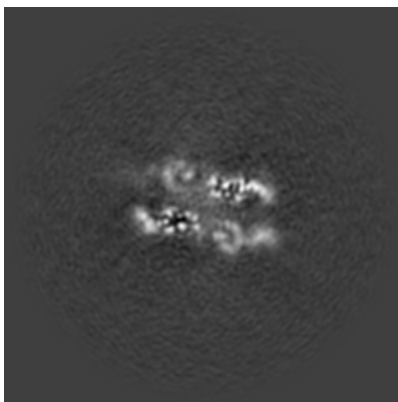
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

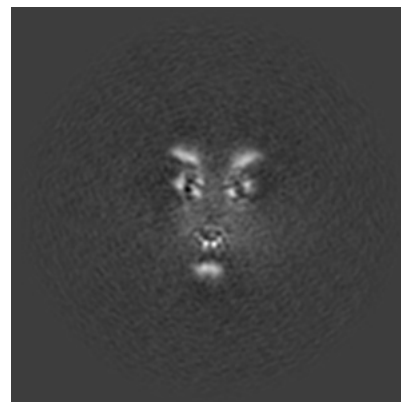
6.2.1 Primary map



X Index: 180

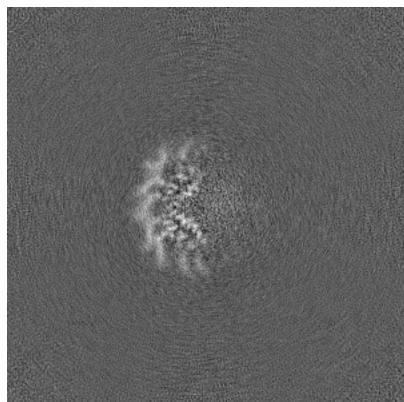


Y Index: 180

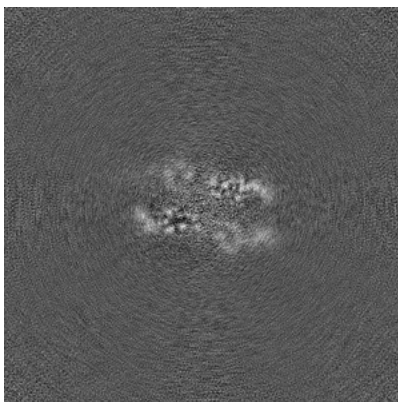


Z Index: 180

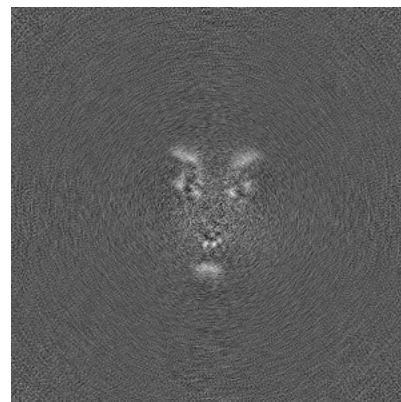
6.2.2 Raw map



X Index: 180



Y Index: 180

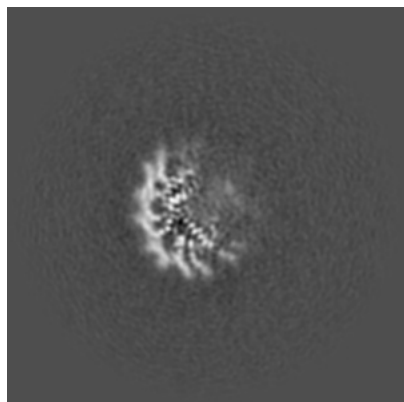


Z Index: 180

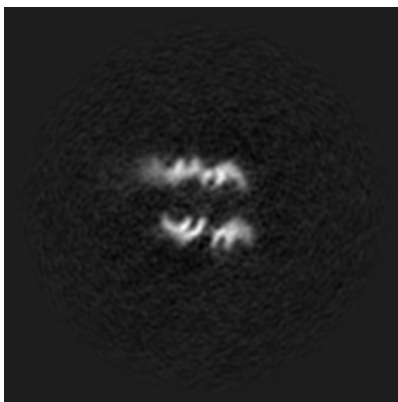
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

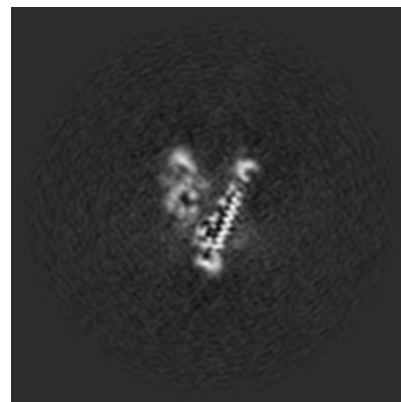
6.3.1 Primary map



X Index: 176

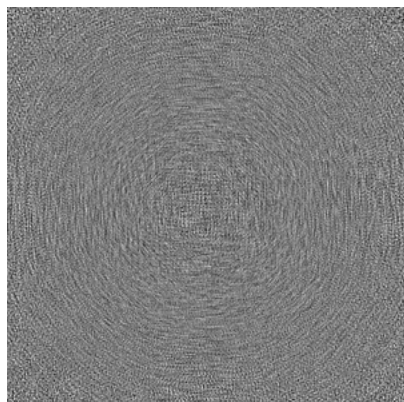


Y Index: 219

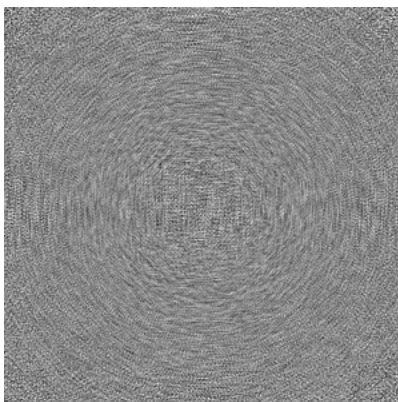


Z Index: 199

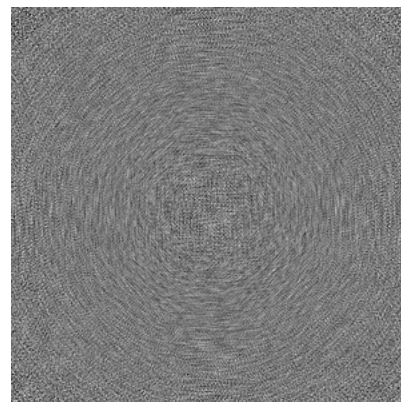
6.3.2 Raw map



X Index: 0



Y Index: 0

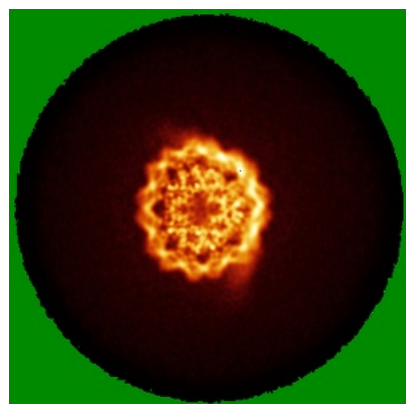


Z Index: 0

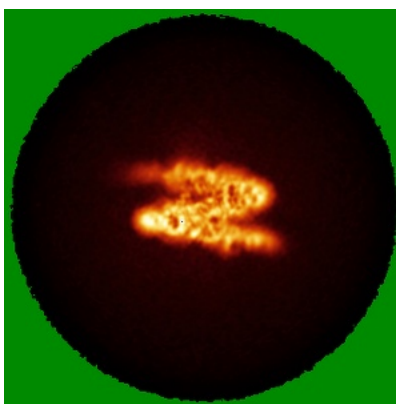
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X

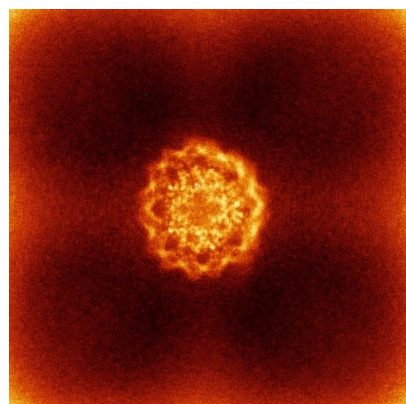


Y

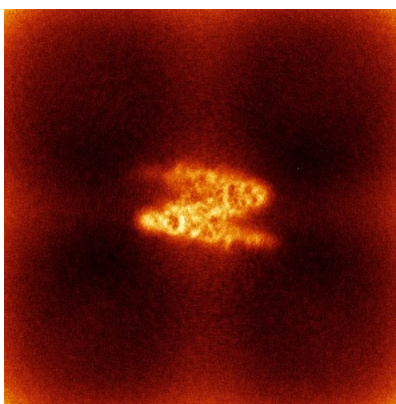


Z

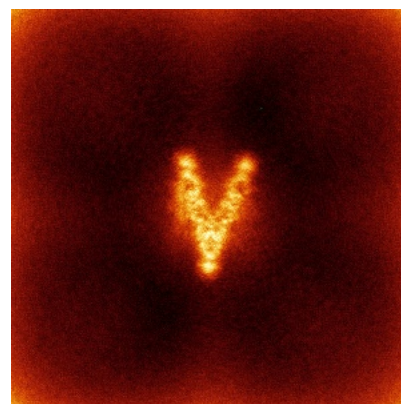
6.4.2 Raw map



X



Y

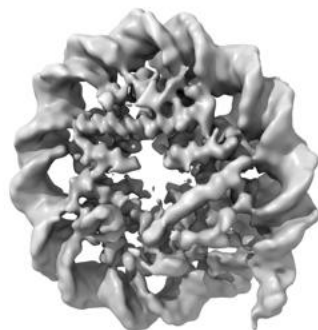


Z

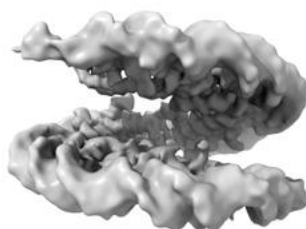
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.108. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

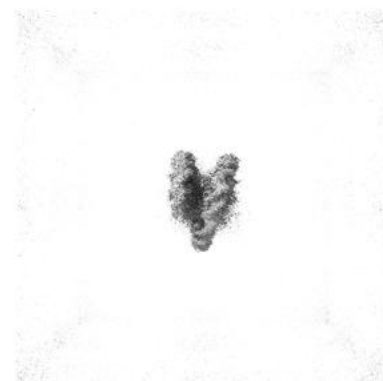
6.5.2 Raw map



X



Y



Z

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.

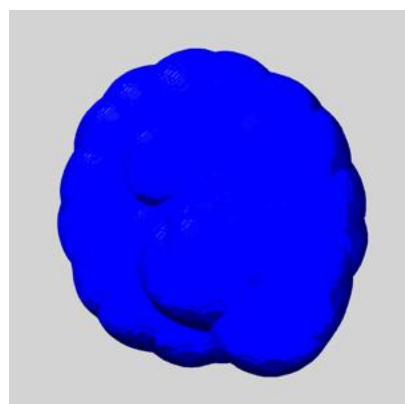
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

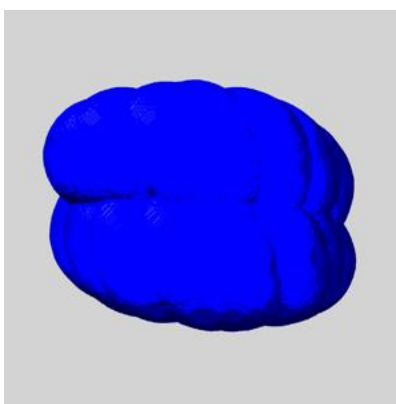
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

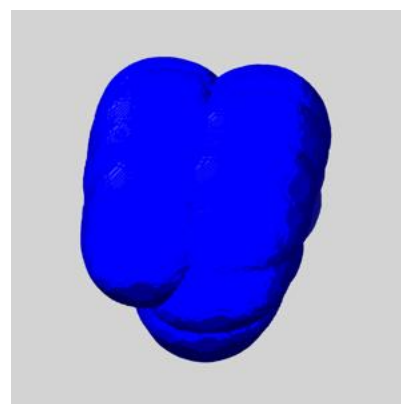
6.6.1 emd_51645_msk_1.map [i](#)



X



Y

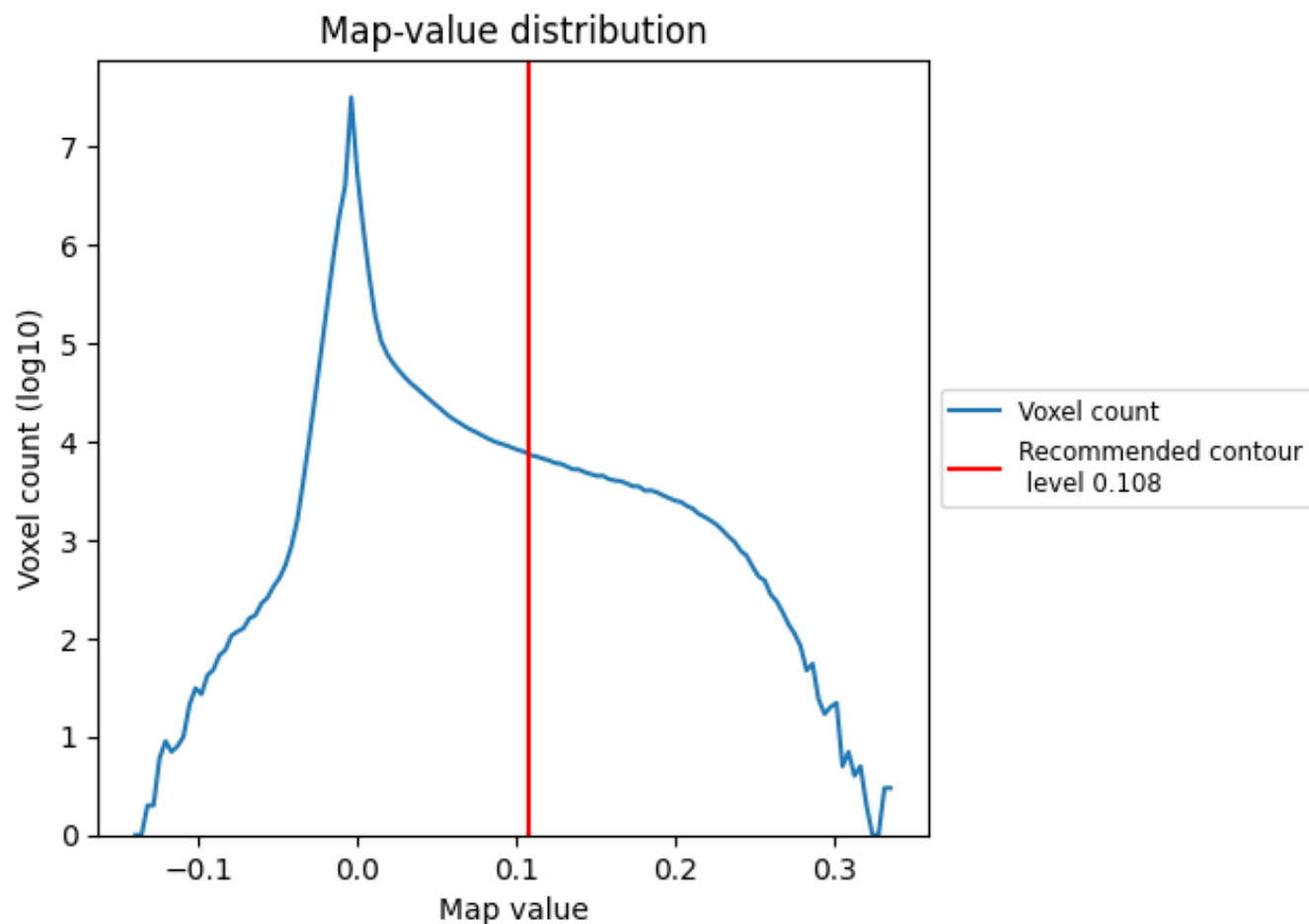


Z

7 Map analysis [i](#)

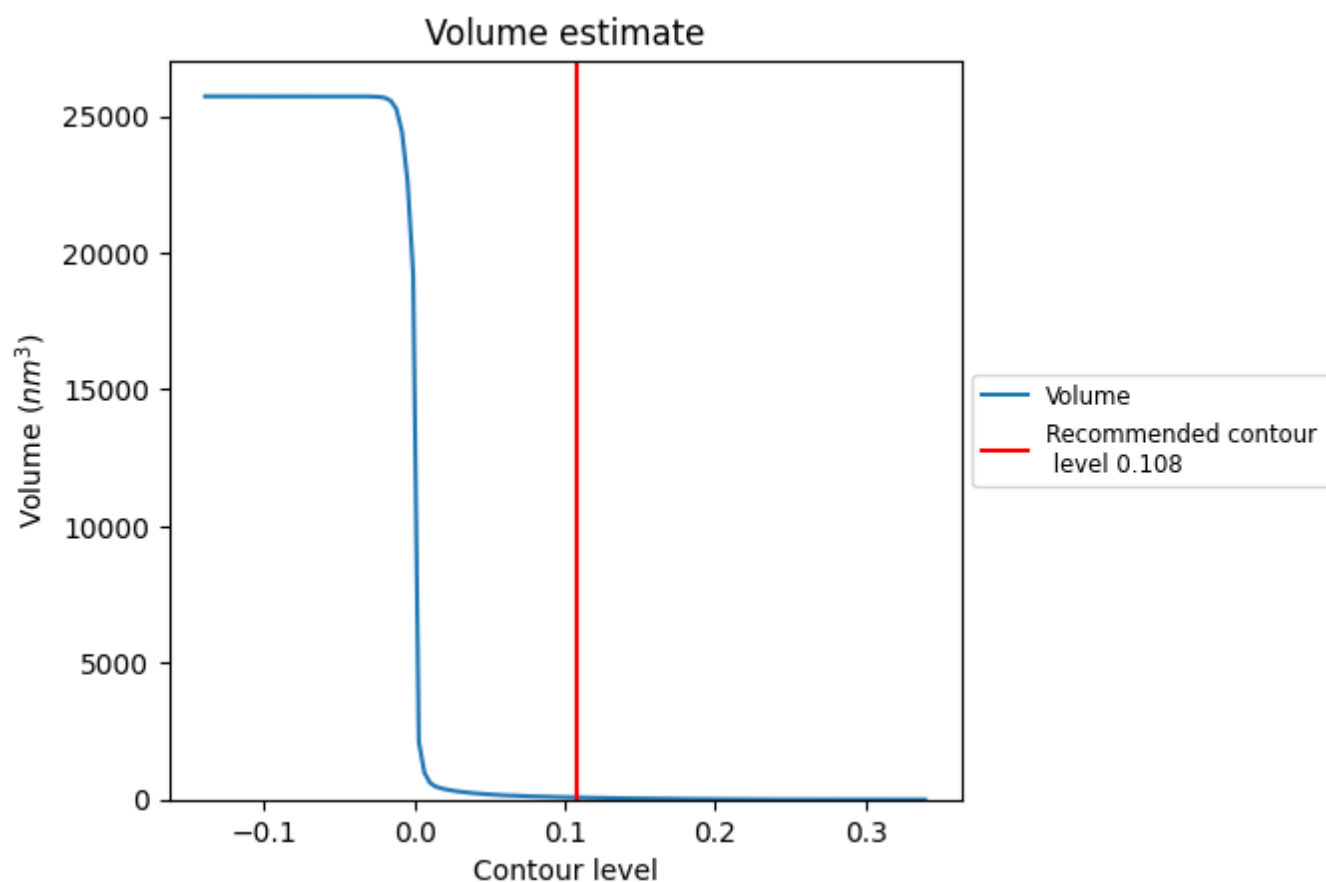
This section contains the results of statistical analysis of the map.

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.

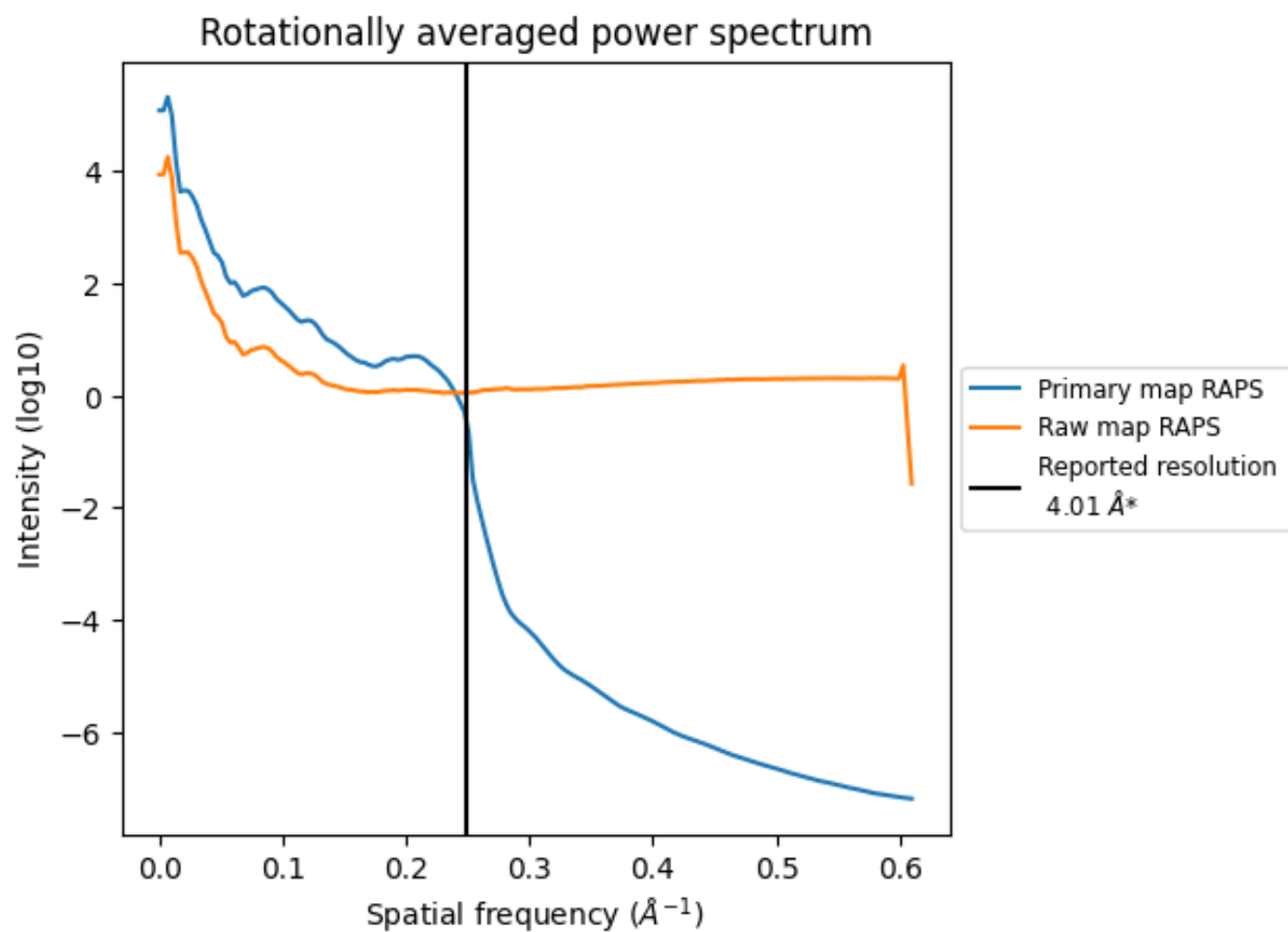
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 76 nm^3 ; this corresponds to an approximate mass of 68 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 ⓘ

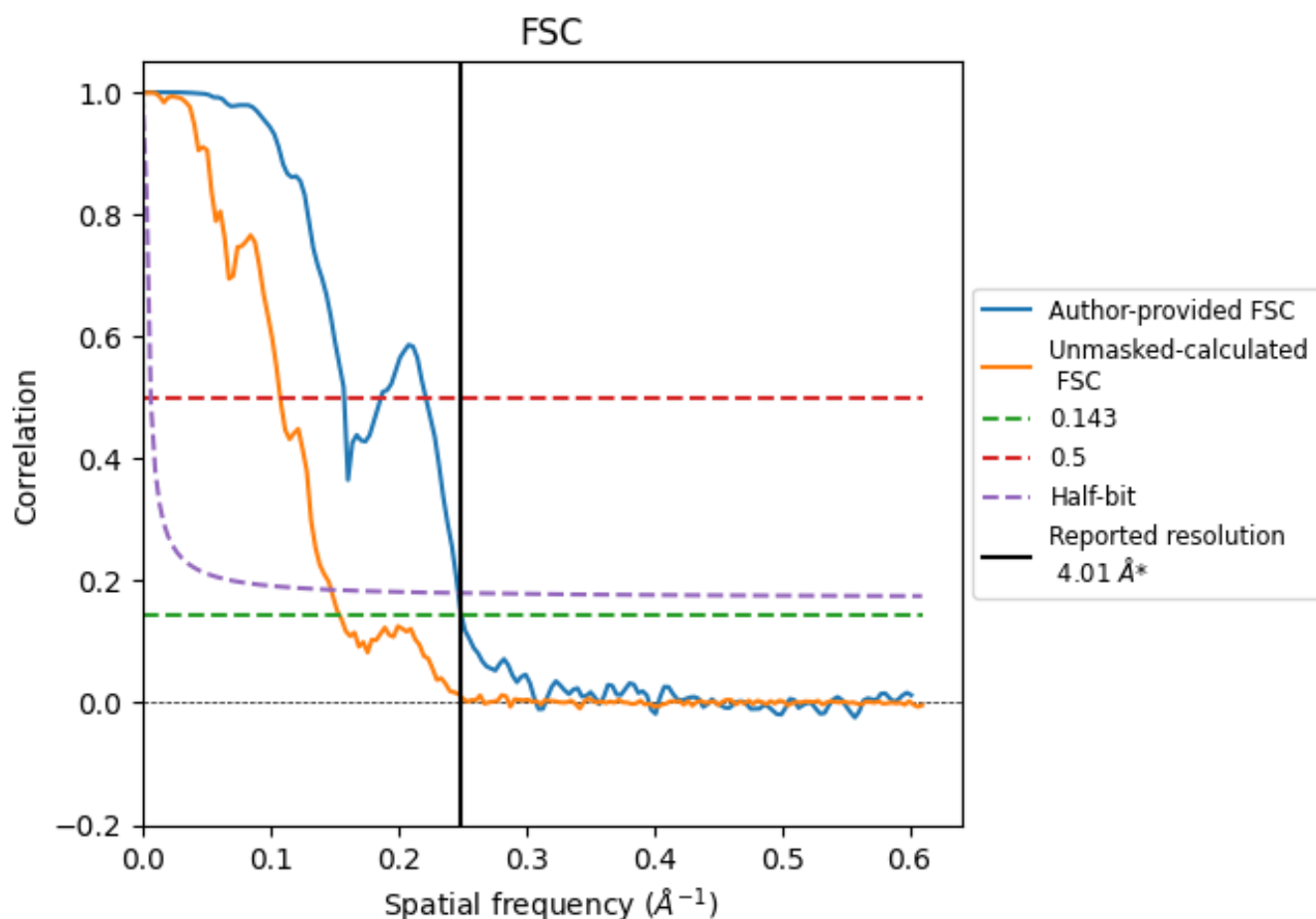


*Reported resolution corresponds to spatial frequency of 0.249 $\text{\AA}^{-1}$

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 0.249 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

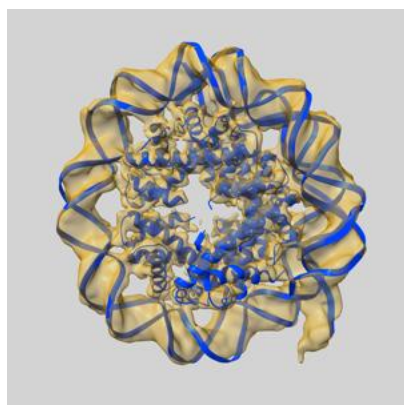
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.01	-	-
Author-provided FSC curve	4.01	6.33	4.05
Unmasked-calculated*	6.48	9.27	6.78

*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 6.48 differs from the reported value 4.01 by more than 10 %

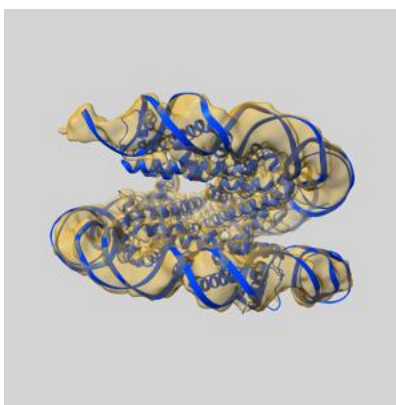
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-51645 and PDB model 9GXA. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

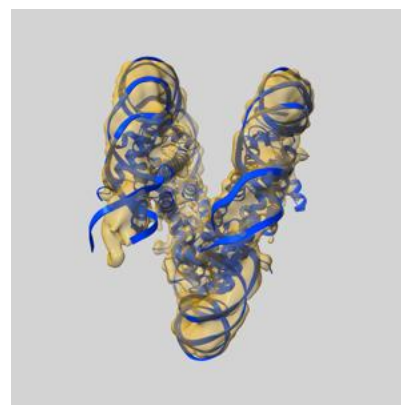
9.1 Map-model overlay [i](#)



X



Y



Z

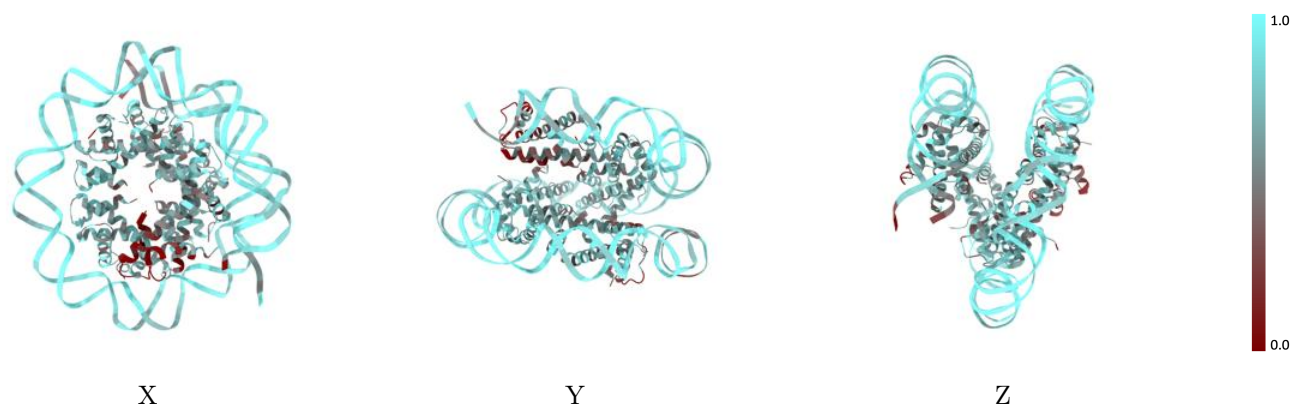
The images above show the 3D surface view of the map at the recommended contour level 0.108 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



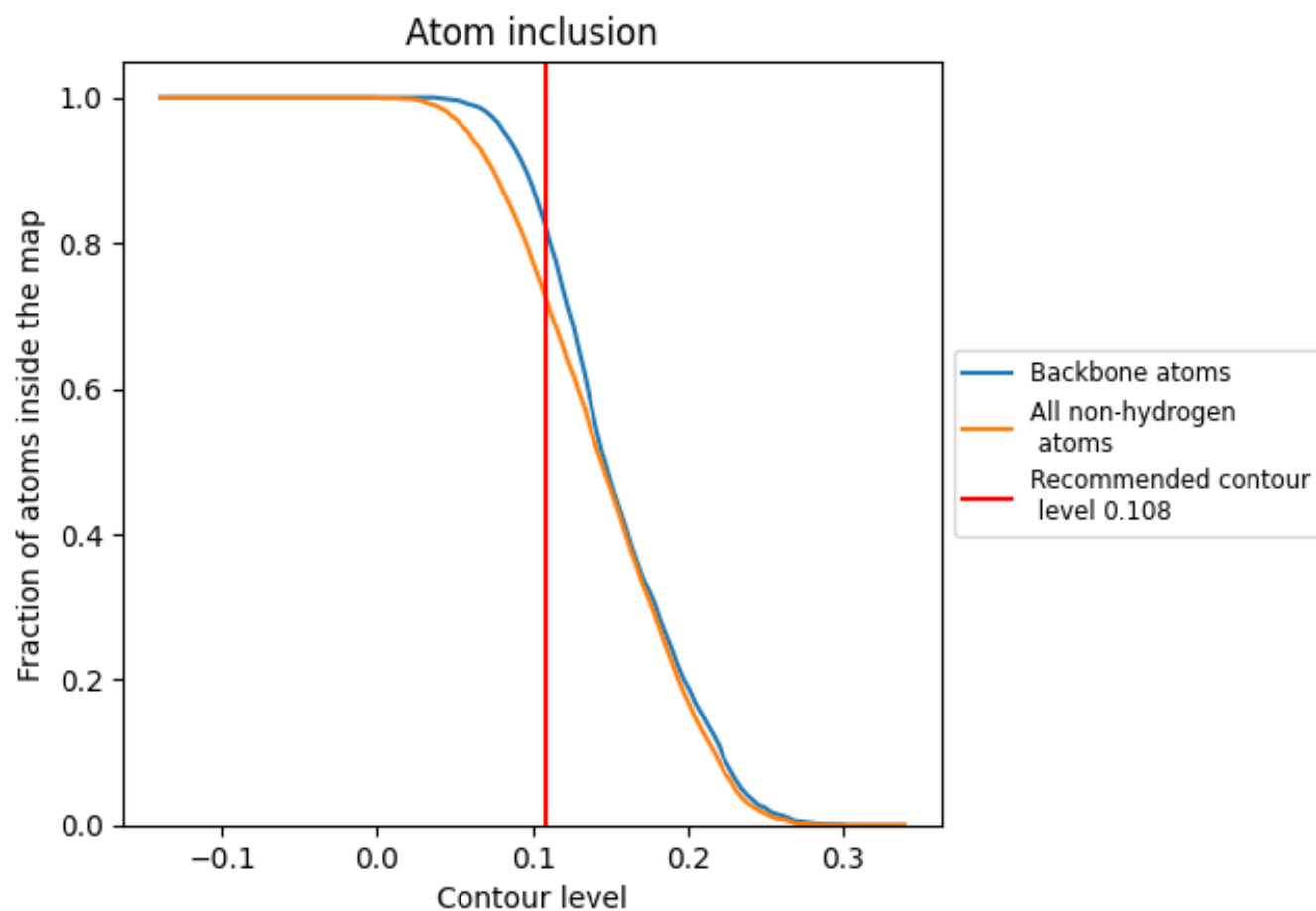
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.108).

9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 73% 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 (0.108) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7280	 0.3180
A	 0.5120	 0.2970
B	 0.4330	 0.2630
C	 0.6660	 0.3850
D	 0.7050	 0.4090
E	 0.7210	 0.4160
F	 0.7080	 0.4010
G	 0.6200	 0.3310
H	 0.6050	 0.3110
I	 0.8310	 0.2800
J	 0.8250	 0.2860

