Full wwPDB EM Validation Report ⓘ

Jan 29, 2025 – 12:11 pm GMT

PDB ID	:	8RUP
EMDB ID	:	EMD-19513
Title	:	Chromosome Passenger Complex (CPC) localization module in complex with H3.T3p-nucleosome
Authors	:	Ruza, R.R.; Barr, F.A.
Deposited on	:	2024-01-31
Resolution	:	2.42 Å(reported)
Based on initial models	:	3AFA, 6YIH

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

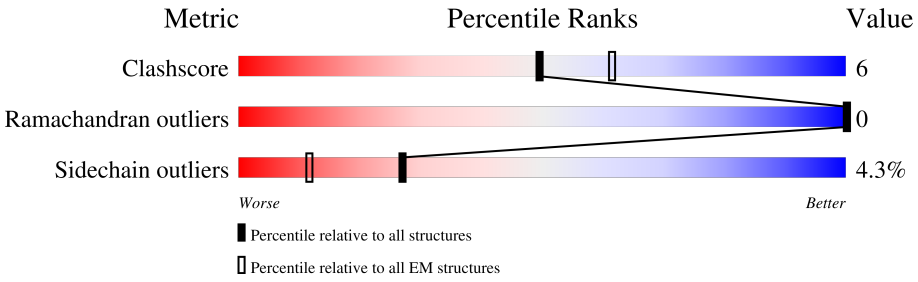
EMDB validation analysis	:	0.0.1.dev113
Mogul	:	1.8.4, CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

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 2.42 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	136	68% · 28%
2	B	103	62% 16% 22%
2	F	103	70% 8% 22%
3	C	130	79% · 17%
3	G	130	79% · 17%
4	D	123	72% 6% 23%
4	H	123	68% 8% · 23%
5	E	135	73% · · 22%

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Mol	Chain	Length	Quality of chain
6	I	152	
7	J	152	
8	K	144	
9	L	76	
10	M	82	

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 25466 atoms, of which 11444 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
1	A	98	Total	C	H	N	O	S	0	0
			1634	506	833	153	139	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	GLY	engineered mutation	UNP P84233

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	80	Total	C	H	N	O	S	0	0
			1303	401	671	122	108	1		
2	F	80	Total	C	H	N	O	S	0	0
			1303	401	671	122	108	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	108	Total	C	H	N	O	0	0
			1676	514	858	162	142		
3	G	108	Total	C	H	N	O	0	0
			1691	517	869	163	142		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	99	ARG	GLY	conflict	UNP P06897
C	123	SER	ALA	conflict	UNP P06897
G	99	ARG	GLY	conflict	UNP P06897
G	123	SER	ALA	conflict	UNP P06897

- Molecule 4 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	95	Total	C	H	N	O	S	0	0
			1472	460	740	131	139	2		
4	H	95	Total	C	H	N	O	S	0	0
			1444	454	723	130	135	2		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	3	MET	-	initiating methionine	UNP P02281
D	32	THR	SER	conflict	UNP P02281
H	3	MET	-	initiating methionine	UNP P02281
H	32	THR	SER	conflict	UNP P02281

- Molecule 5 is a protein called Histone H3.

Mol	Chain	Residues	Atoms							AltConf	Trace
5	E	105	Total	C	H	N	O	P	S	0	0
			1723	532	875	164	148	1	3		

- Molecule 6 is a DNA chain called DNA (147-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
6	I	147	Total	C	H	N	O	P	0	0
			4679	1434	1648	570	880	147		

- Molecule 7 is a DNA chain called DNA (147-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
7	J	147	Total	C	H	N	O	P	0	0
			4643	1423	1650	542	882	146		

- Molecule 8 is a protein called Baculoviral IAP repeat-containing protein 5.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	K	136	Total	C	H	N	O	S	0	0
			2031	675	976	178	195	7		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	-1	GLY	-	expression tag	UNP O15392
K	0	PRO	-	expression tag	UNP O15392

- Molecule 9 is a protein called Borealin.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	L	75	Total	C	H	N	O	S	0	0
			1264	396	639	117	111	1		

- Molecule 10 is a protein called Inner centromere protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	M	41	Total	C	H	N	O	S	0	0
			602	202	291	49	55	5		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	-1	GLY	-	expression tag	UNP Q9NQS7
M	0	PRO	-	expression tag	UNP Q9NQS7

- Molecule 11 is ZINC ION (three-letter code: ZN) (formula: Zn).

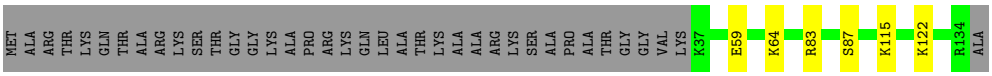
Mol	Chain	Residues	Atoms		AltConf
11	K	1	Total	Zn	0
			1	1	

3 Residue-property plots (i)

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.

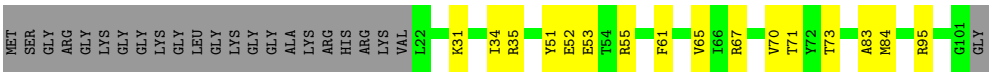
• Molecule 1: Histone H3.2

Chain A:



• Molecule 2: Histone H4

Chain B:



• Molecule 2: Histone H4

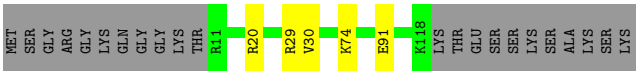
Chain F:

• Molecule 3: Histone H2A type 1

Chain C:

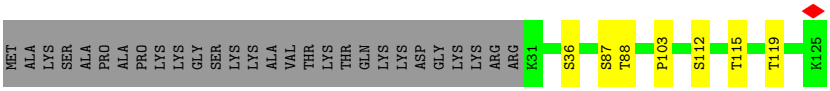
• Molecule 3: Histone H2A type 1

Chain G:



• Molecule 4: Histone H2B 1.1

Chain D:

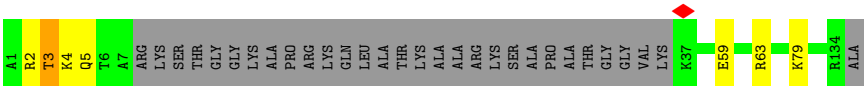


• Molecule 4: Histone H2B 1.1

Chain H:

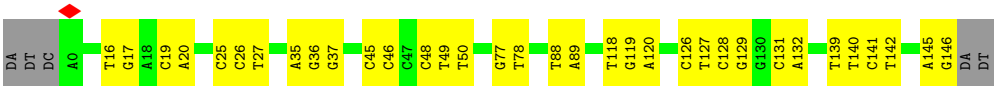
• Molecule 5: Histone H3

Chain E:



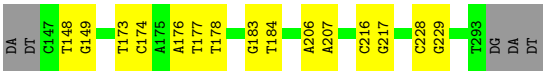
• Molecule 6: DNA (147-MER)

Chain I:



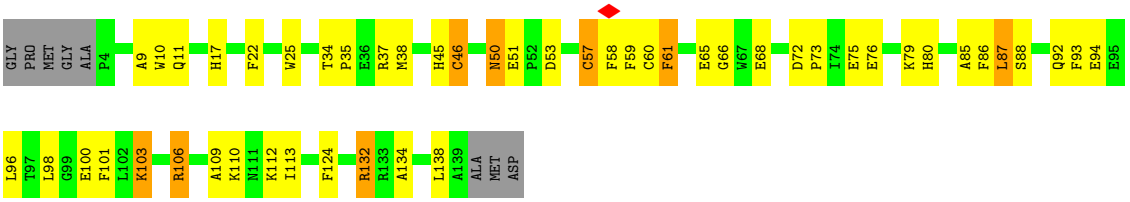
• Molecule 7: DNA (147-MER)

Chain J:



• Molecule 8: Baculoviral IAP repeat-containing protein 5

Chain K:

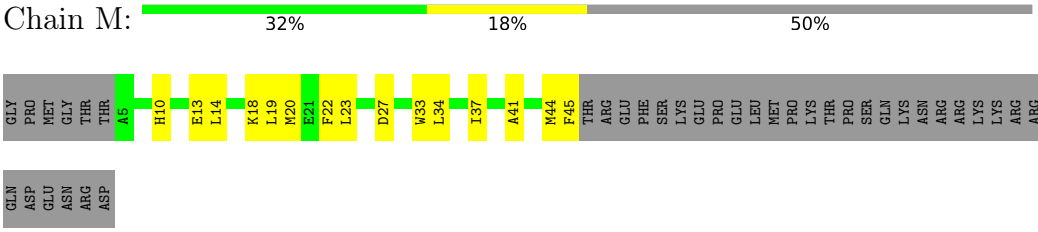


• Molecule 9: Borealin

Chain L:

• Molecule 10: Inner centromere protein

Chain M:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	119063	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; CTF amplitude correction first performed following movie motion correction, and CTF parameters were later refined during the refinement of the final reconstruction	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	43.86	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2300	Depositor
Magnification	58009	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.911	Depositor
Minimum map value	-0.000	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	318.72, 318.72, 318.72	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.34	0/813	0.61	0/1093
2	B	0.34	0/639	0.62	0/857
2	F	0.34	0/639	0.60	0/857
3	C	0.33	0/828	0.57	0/1121
3	G	0.31	0/832	0.58	0/1125
4	D	0.34	0/743	0.50	0/1003
4	H	0.33	0/732	0.51	0/990
5	E	0.33	0/847	0.58	0/1136
6	I	0.66	0/3404	0.91	0/5256
7	J	0.65	0/3353	0.91	0/5169
8	K	0.33	0/1085	0.50	0/1470
9	L	0.35	0/633	0.61	0/846
10	M	0.32	0/317	0.55	0/430
All	All	0.50	0/14865	0.75	0/21353

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

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	801	833	831	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	632	671	671	7	0
2	F	632	671	671	4	0
3	C	818	858	858	3	0
3	G	822	869	869	3	0
4	D	732	740	740	3	0
4	H	721	723	723	5	0
5	E	848	875	875	5	0
6	I	3031	1648	1649	28	0
7	J	2993	1650	1652	8	0
8	K	1055	976	974	50	0
9	L	625	639	650	27	0
10	M	311	291	289	13	0
11	K	1	0	0	0	0
All	All	14022	11444	11452	140	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 6.

All (140) 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)
9:L:24:PHE:HE2	10:M:41:ALA:HB2	1.50	0.76
10:M:23:LEU:HD13	10:M:27:ASP:OD2	1.85	0.76
2:B:52:GLU:OE2	2:B:55:ARG:NH1	2.21	0.73
3:C:42:ARG:NH1	4:D:88:THR:OG1	2.22	0.72
3:G:91:GLU:N	3:G:91:GLU:OE1	2.24	0.71
8:K:38:MET:SD	8:K:38:MET:N	2.66	0.69
9:L:42:ASP:HB3	10:M:22:PHE:HE1	1.59	0.67
5:E:59:GLU:OE1	5:E:59:GLU:N	2.27	0.67
9:L:24:PHE:CE2	10:M:41:ALA:HB2	2.30	0.67
8:K:132:ARG:HH12	9:L:25:LEU:HD22	1.61	0.64
1:A:59:GLU:OE1	1:A:59:GLU:N	2.29	0.64
2:F:26:ILE:HD13	2:F:59:LYS:HD3	1.79	0.64
4:H:76:GLU:OE1	4:H:79:ARG:NH1	2.35	0.59
5:E:2:ARG:NH2	8:K:65:GLU:OE2	2.34	0.59
9:L:71:LEU:O	9:L:75:ALA:N	2.34	0.59
8:K:110:LYS:HE3	8:K:110:LYS:HA	1.85	0.58
2:F:82:THR:HG22	2:F:84:MET:H	1.70	0.57
6:I:118:DT:H2"	6:I:119:DG:C8	2.39	0.57
8:K:72:ASP:O	8:K:72:ASP:OD1	2.23	0.57
8:K:134:ALA:O	8:K:138:LEU:HG	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:K:53:ASP:N	8:K:53:ASP:OD1	2.36	0.56
9:L:56:ILE:CD1	9:L:59:LEU:HD11	2.35	0.56
8:K:87:LEU:HD12	8:K:88:SER:N	2.22	0.55
10:M:18:LYS:N	10:M:18:LYS:HD3	2.22	0.55
8:K:72:ASP:OD1	8:K:76:GLU:OE2	2.26	0.53
2:B:70:VAL:HA	2:B:73:THR:HG22	1.90	0.53
8:K:46:CYS:O	8:K:46:CYS:SG	2.66	0.53
6:I:139:DT:H2'	6:I:140:DT:H72	1.89	0.52
10:M:10:HIS:O	10:M:14:LEU:HD13	2.09	0.52
9:L:63:LYS:HA	9:L:66:ARG:HB3	1.91	0.52
8:K:92:GLN:OE1	8:K:93:PHE:N	2.38	0.52
10:M:33:TRP:CH2	10:M:37:ILE:HD13	2.45	0.52
8:K:132:ARG:NH1	9:L:25:LEU:HD22	2.25	0.52
6:I:126:DC:H2'	6:I:127:DT:H72	1.91	0.52
7:J:173:DT:H2"	7:J:174:DC:C6	2.45	0.51
10:M:33:TRP:O	10:M:37:ILE:HG23	2.10	0.51
8:K:103:LYS:HA	8:K:103:LYS:HE2	1.92	0.51
9:L:38:GLN:NE2	9:L:42:ASP:OD2	2.43	0.51
8:K:34:THR:O	8:K:38:MET:SD	2.69	0.51
8:K:86:PHE:HE1	8:K:93:PHE:CE2	2.28	0.51
7:J:177:DT:C6	7:J:178:DT:H72	2.46	0.50
6:I:48:DC:H2'	6:I:49:DT:H72	1.93	0.50
6:I:88:DT:H2"	6:I:89:DA:C8	2.47	0.50
2:B:61:PHE:O	2:B:65:VAL:HG23	2.12	0.50
8:K:109:ALA:O	8:K:113:ILE:HD13	2.12	0.50
8:K:61:PHE:CD1	8:K:61:PHE:N	2.77	0.49
4:D:115:THR:O	4:D:119:THR:HG22	2.12	0.49
4:H:33:ARG:HB3	4:H:33:ARG:NH1	2.26	0.49
9:L:43:ARG:O	9:L:47:LEU:HG	2.13	0.49
6:I:77:DG:H2"	6:I:78:DT:H71	1.95	0.48
9:L:62:PRO:O	9:L:66:ARG:N	2.42	0.48
9:L:56:ILE:HD12	9:L:59:LEU:HD11	1.94	0.48
8:K:65:GLU:HG3	8:K:66:GLY:N	2.29	0.48
9:L:70:TRP:CG	9:L:71:LEU:N	2.81	0.48
8:K:9:ALA:HB3	8:K:10:TRP:CZ3	2.49	0.48
8:K:106:ARG:NE	8:K:110:LYS:HD3	2.29	0.48
10:M:19:LEU:HD23	10:M:20:MET:N	2.29	0.48
8:K:22:PHE:CD1	8:K:45:HIS:HB2	2.49	0.47
6:I:36:DG:C2'	6:I:37:DG:C8	2.97	0.47
6:I:49:DT:H2'	6:I:50:DT:H71	1.95	0.47
8:K:79:LYS:HG3	8:K:80:HIS:ND1	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:L:63:LYS:O	9:L:66:ARG:N	2.47	0.47
9:L:63:LYS:O	9:L:67:GLU:OE1	2.32	0.47
9:L:63:LYS:HG3	9:L:64:ALA:N	2.30	0.47
8:K:22:PHE:HE2	8:K:35:PRO:O	1.98	0.47
8:K:72:ASP:CG	8:K:75:GLU:OE1	2.53	0.47
9:L:53:LEU:O	9:L:56:ILE:HG22	2.15	0.46
8:K:68:GLU:HA	8:K:68:GLU:OE1	2.16	0.46
3:G:30:VAL:HG13	4:H:70:PHE:HE2	1.80	0.46
1:A:87:SER:OG	2:B:83:ALA:HB2	2.16	0.46
8:K:94:GLU:N	8:K:94:GLU:CD	2.69	0.46
6:I:26:DC:H2"	6:I:27:DT:C5	2.50	0.46
7:J:216:DC:H2"	7:J:217:DG:C8	2.51	0.46
6:I:36:DG:N2	6:I:37:DG:C2	2.84	0.45
6:I:45:DC:H2"	6:I:46:DC:C6	2.52	0.45
6:I:141:DC:C2'	6:I:142:DT:H72	2.47	0.45
6:I:36:DG:H2"	6:I:37:DG:O5'	2.17	0.45
3:G:20:ARG:HD2	4:H:121:TYR:CE1	2.51	0.45
8:K:9:ALA:HB3	8:K:10:TRP:CE3	2.52	0.45
8:K:96:LEU:HB2	9:L:70:TRP:CZ3	2.51	0.45
9:L:44:GLN:OE1	9:L:47:LEU:HD12	2.17	0.45
5:E:3:TPO:HG23	5:E:4:LYS:N	2.32	0.45
8:K:79:LYS:O	8:K:79:LYS:HD2	2.17	0.45
8:K:86:PHE:CE1	8:K:93:PHE:CZ	3.06	0.44
6:I:36:DG:H2'	6:I:37:DG:C8	2.52	0.44
8:K:58:PHE:O	8:K:61:PHE:HE1	2.00	0.44
8:K:25:TRP:CH2	8:K:38:MET:HG3	2.53	0.44
8:K:35:PRO:HA	8:K:38:MET:HG2	2.00	0.44
8:K:11:GLN:HB3	8:K:17:HIS:CD2	2.52	0.44
9:L:44:GLN:OE1	9:L:44:GLN:HA	2.18	0.44
9:L:63:LYS:O	9:L:66:ARG:HB3	2.17	0.44
9:L:62:PRO:HG2	9:L:65:LEU:HB3	1.99	0.44
6:I:36:DG:H2"	6:I:37:DG:C8	2.52	0.44
7:J:228:DC:H2"	7:J:229:DG:C8	2.53	0.44
6:I:25:DC:H2"	6:I:26:DC:C6	2.53	0.44
7:J:206:DA:H2"	7:J:207:DA:C8	2.52	0.44
8:K:57:CYS:SG	8:K:59:PHE:N	2.82	0.44
6:I:16:DT:H2"	6:I:17:DG:C8	2.53	0.43
6:I:35:DA:H2"	6:I:36:DG:C8	2.53	0.43
6:I:36:DG:H4'	6:I:37:DG:OP1	2.17	0.43
8:K:37:ARG:HB3	8:K:73:PRO:HG2	2.01	0.43
2:F:84:MET:HA	2:F:84:MET:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:K:57:CYS:O	8:K:61:PHE:HA	2.18	0.43
9:L:69:ASN:OD1	9:L:71:LEU:N	2.52	0.43
5:E:3:TPO:O	5:E:5:GLN:NE2	2.51	0.42
4:H:48:VAL:HG23	4:H:49:HIS:N	2.34	0.42
6:I:128:DC:H2"	6:I:129:DG:N7	2.34	0.42
8:K:94:GLU:HA	9:L:70:TRP:HH2	1.84	0.42
6:I:36:DG:C2	6:I:37:DG:C2	3.07	0.42
7:J:183:DG:H2"	7:J:184:DT:C6	2.55	0.42
6:I:119:DG:H1'	6:I:120:DA:N7	2.35	0.42
3:C:92:GLU:OE1	4:D:103:PRO:O	2.38	0.42
6:I:77:DG:H2"	6:I:78:DT:C7	2.50	0.42
7:J:176:DA:C2'	7:J:177:DT:H71	2.50	0.42
8:K:85:ALA:O	8:K:88:SER:N	2.47	0.42
8:K:86:PHE:CE1	8:K:93:PHE:CE2	3.07	0.42
8:K:106:ARG:CZ	8:K:110:LYS:HD3	2.50	0.42
2:B:31:LYS:O	2:B:35:ARG:HG2	2.19	0.41
2:B:34:ILE:HG21	2:B:51:TYR:HD1	1.85	0.41
3:C:71:ARG:HH11	3:C:71:ARG:HG2	1.84	0.41
5:E:2:ARG:N	8:K:65:GLU:O	2.42	0.41
6:I:139:DT:C2'	6:I:140:DT:H72	2.50	0.41
8:K:96:LEU:HD22	8:K:100:GLU:HB2	2.01	0.41
8:K:61:PHE:HE2	8:K:112:LYS:N	2.18	0.41
8:K:98:LEU:HD21	9:L:73:TYR:CD2	2.55	0.41
2:B:67:ARG:O	2:B:71:THR:HG22	2.21	0.41
8:K:50:ASN:HD22	8:K:51:GLU:N	2.19	0.41
8:K:57:CYS:SG	8:K:60:CYS:N	2.80	0.41
9:L:31:GLU:CD	10:M:33:TRP:CZ2	2.94	0.41
6:I:48:DC:H2"	6:I:49:DT:C6	2.56	0.41
8:K:124:PHE:CE1	10:M:34:LEU:HD11	2.56	0.41
8:K:22:PHE:HD2	8:K:35:PRO:HB3	1.85	0.41
6:I:19:DC:H2"	6:I:20:DA:C8	2.56	0.40
6:I:131:DC:H2"	6:I:132:DA:H5'	2.04	0.40
10:M:44:MET:HE1	10:M:45:PHE:CE2	2.56	0.40
6:I:145:DA:H2"	6:I:146:DG:C8	2.57	0.40
7:J:148:DT:H2"	7:J:149:DG:C8	2.57	0.40
2:F:35:ARG:O	2:F:39:ARG:HG2	2.21	0.40
8:K:98:LEU:HD12	9:L:65:LEU:O	2.22	0.40
10:M:10:HIS:O	10:M:13:GLU:OE1	2.39	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	96/136 (71%)	95 (99%)	1 (1%)	0	100	100
2	B	78/103 (76%)	78 (100%)	0	0	100	100
2	F	78/103 (76%)	77 (99%)	1 (1%)	0	100	100
3	C	106/130 (82%)	102 (96%)	4 (4%)	0	100	100
3	G	106/130 (82%)	103 (97%)	3 (3%)	0	100	100
4	D	93/123 (76%)	92 (99%)	1 (1%)	0	100	100
4	H	93/123 (76%)	93 (100%)	0	0	100	100
5	E	100/135 (74%)	99 (99%)	1 (1%)	0	100	100
8	K	134/144 (93%)	132 (98%)	2 (2%)	0	100	100
9	L	73/76 (96%)	69 (94%)	4 (6%)	0	100	100
10	M	39/82 (48%)	36 (92%)	3 (8%)	0	100	100
All	All	996/1285 (78%)	976 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	84/111 (76%)	80 (95%)	4 (5%)	21	35
2	B	64/79 (81%)	61 (95%)	3 (5%)	22	36
2	F	64/79 (81%)	62 (97%)	2 (3%)	35	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	C	81/102 (79%)	79 (98%)	2 (2%)	42 61
3	G	82/102 (80%)	80 (98%)	2 (2%)	44 63
4	D	78/103 (76%)	75 (96%)	3 (4%)	28 45
4	H	75/103 (73%)	70 (93%)	5 (7%)	13 21
5	E	85/109 (78%)	83 (98%)	2 (2%)	44 63
8	K	106/124 (86%)	97 (92%)	9 (8%)	8 13
9	L	67/70 (96%)	64 (96%)	3 (4%)	23 38
10	M	31/77 (40%)	31 (100%)	0	100 100
All	All	817/1059 (77%)	782 (96%)	35 (4%)	27 40

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LYS
1	A	83	ARG
1	A	115	LYS
1	A	122	LYS
2	B	53	GLU
2	B	84	MET
2	B	95	ARG
3	C	36	LYS
3	C	113	SER
4	D	36	SER
4	D	87	SER
4	D	112	SER
5	E	63	ARG
5	E	79	LYS
2	F	47	SER
2	F	95	ARG
3	G	29	ARG
3	G	74	LYS
4	H	33	ARG
4	H	47	GLN
4	H	56	SER
4	H	64	SER
4	H	116	LYS
8	K	46	CYS
8	K	50	ASN
8	K	57	CYS

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Mol	Chain	Res	Type
8	K	61	PHE
8	K	87	LEU
8	K	101	PHE
8	K	103	LYS
8	K	106	ARG
8	K	132	ARG
9	L	8	SER
9	L	66	ARG
9	L	73	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
8	K	50	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	TPO	E	3	5	8,10,11	0.88	0	10,14,16	1.48	1 (10%)

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
5	TPO	E	3	5	-	1/9/11/13	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	E	3	TPO	P-OG1-CB	-4.10	110.83	123.21

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	3	TPO	O-C-CA-CB

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	3	TPO	2	0

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

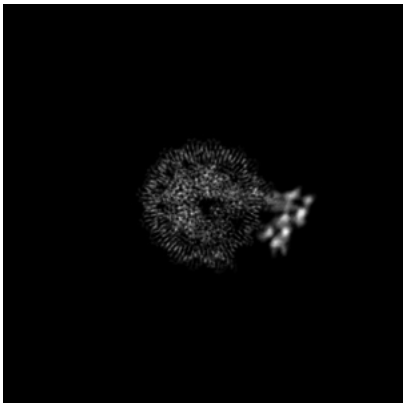
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-19513. These allow visual inspection of the internal detail of the map and identification of artifacts.

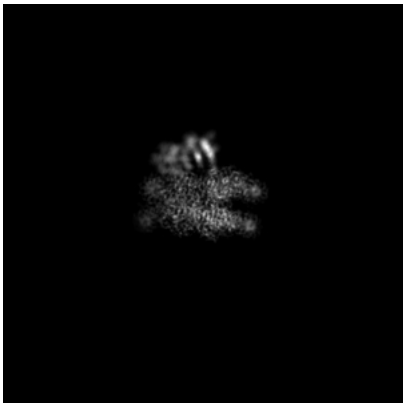
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

6.1.1 Primary map



X



Y

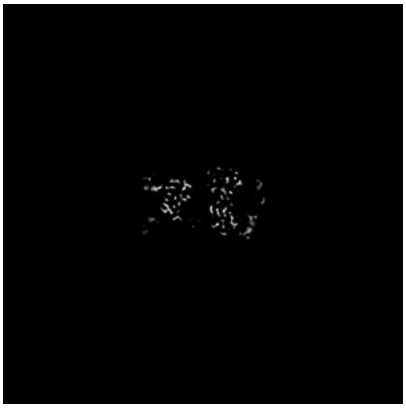
Z

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

6.2 Central slices

6.2.1 Primary map

X Index: 192



Y Index: 192

Z Index: 192

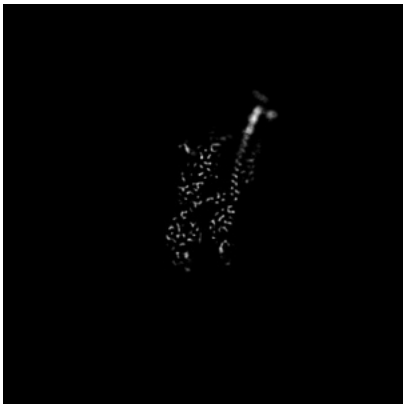
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

6.3.1 Primary map

X Index: 245

Y Index: 271

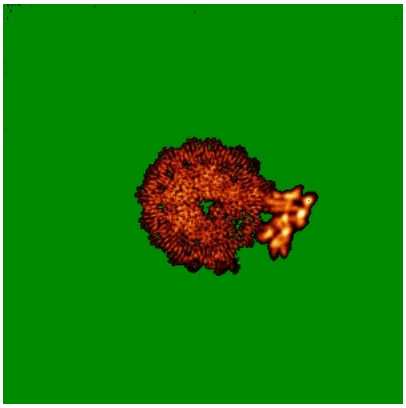


Z Index: 200

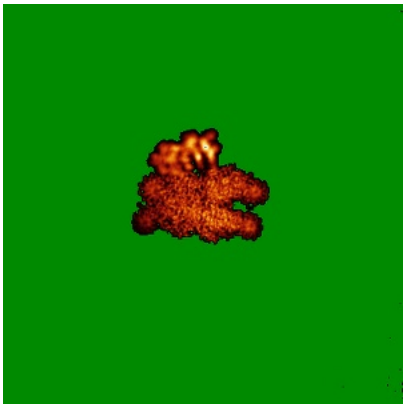
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



X



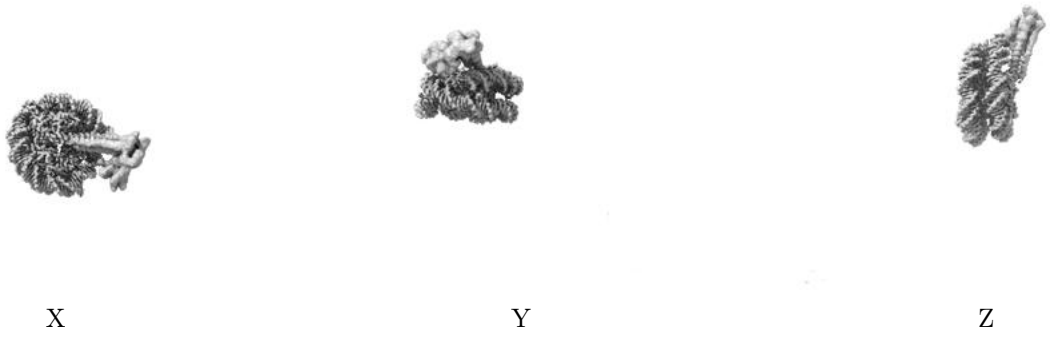
Y

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 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

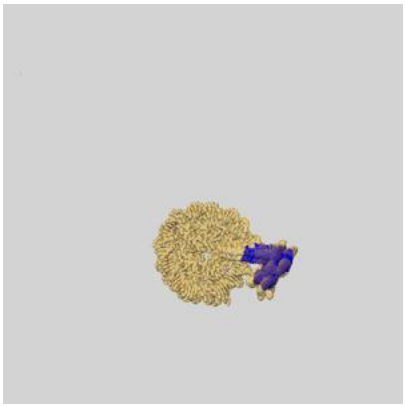
6.6 Mask visualisation (i)

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

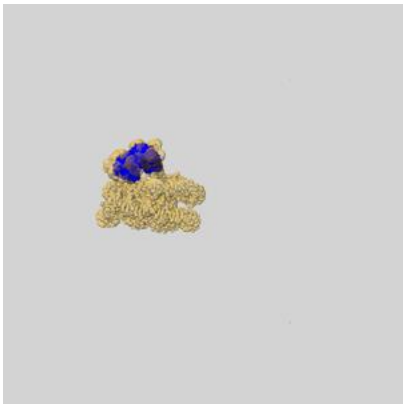
A mask typically either:

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

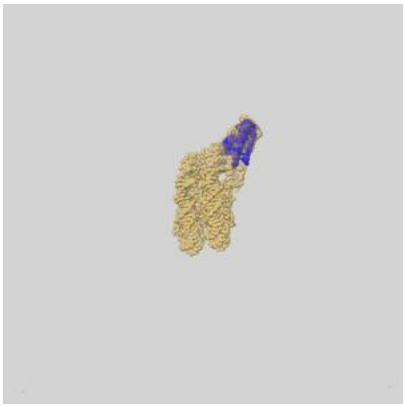
6.6.1 emd_19513_msk_1.map (i)



X



Y

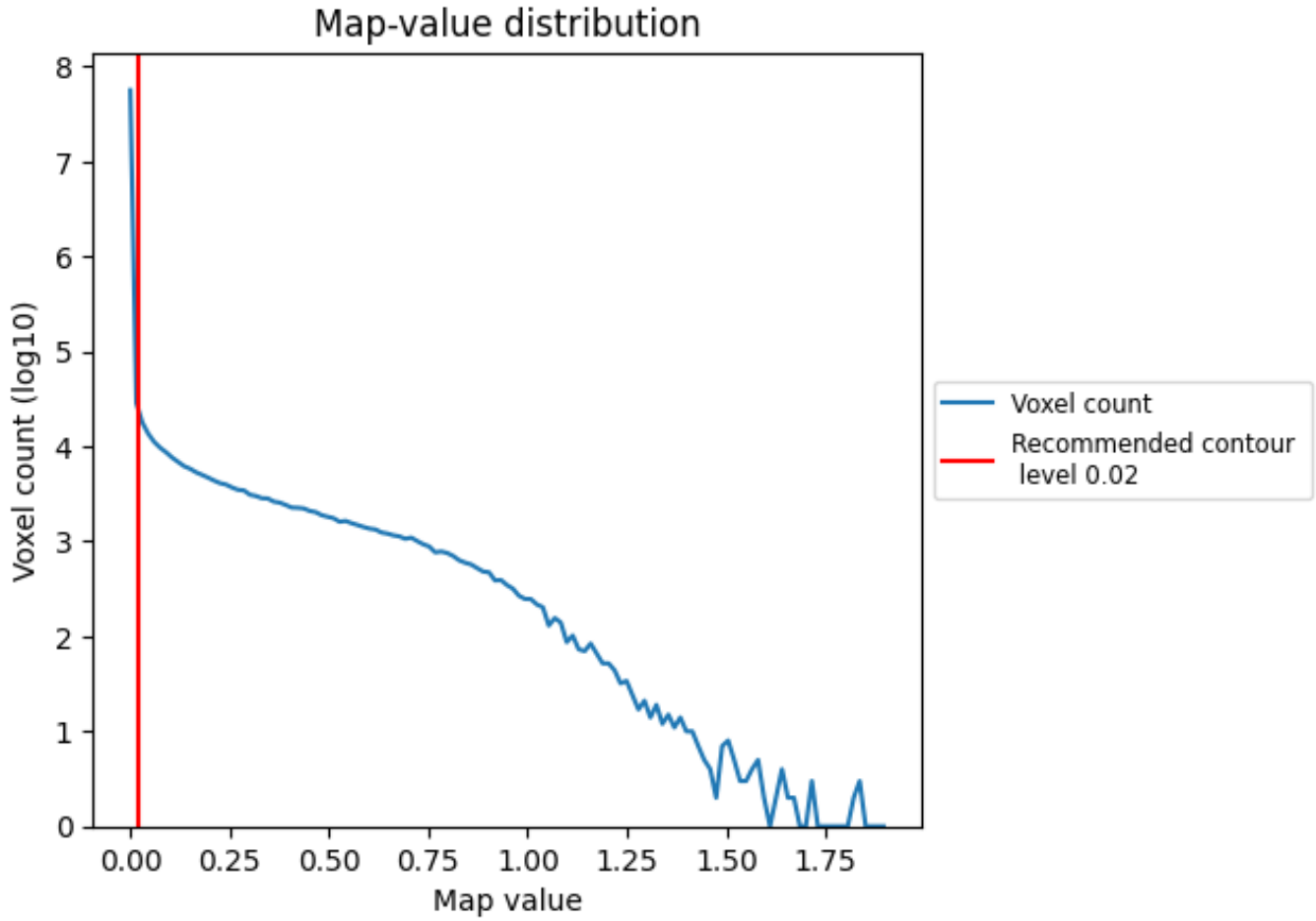


Z

7 Map analysis

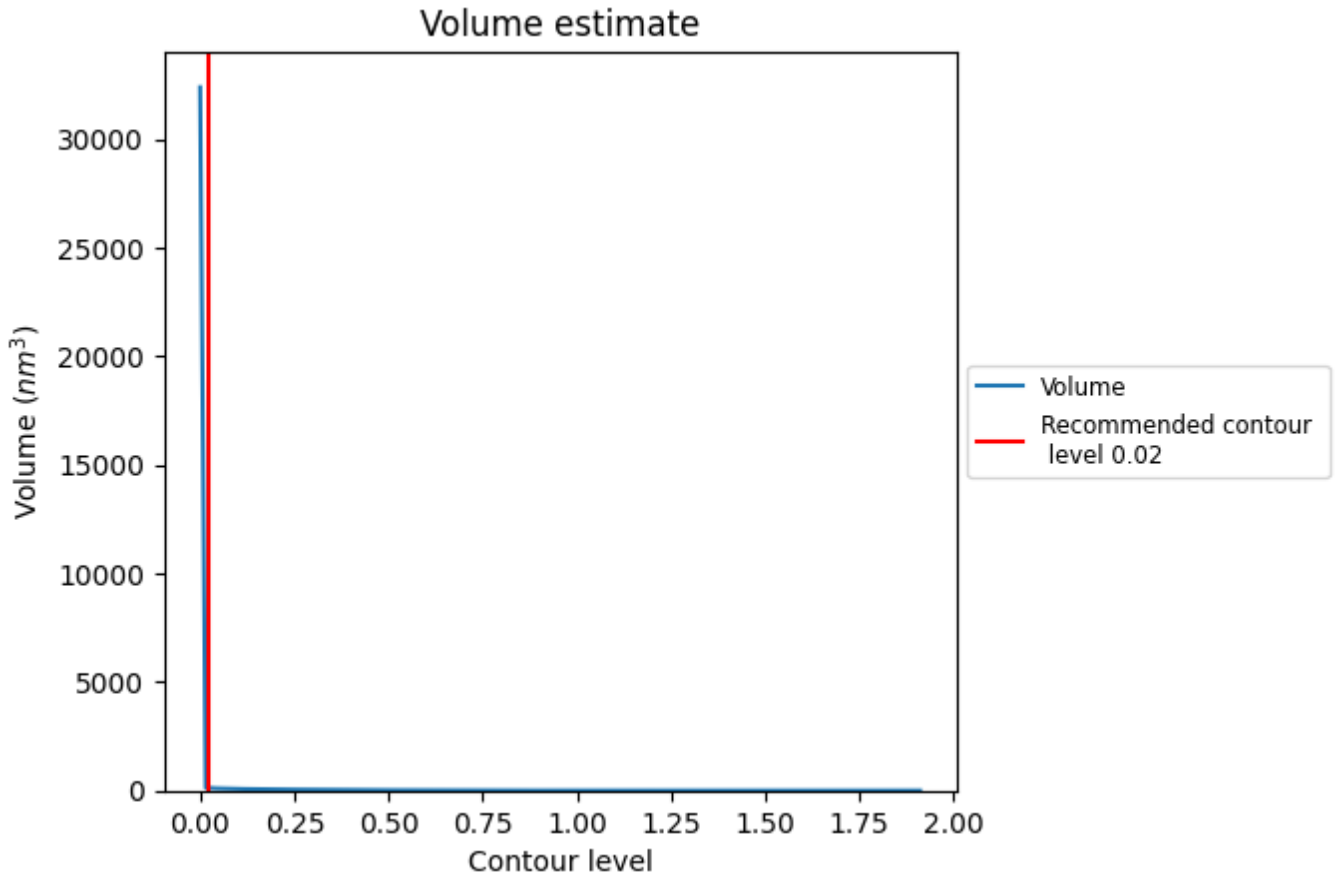
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

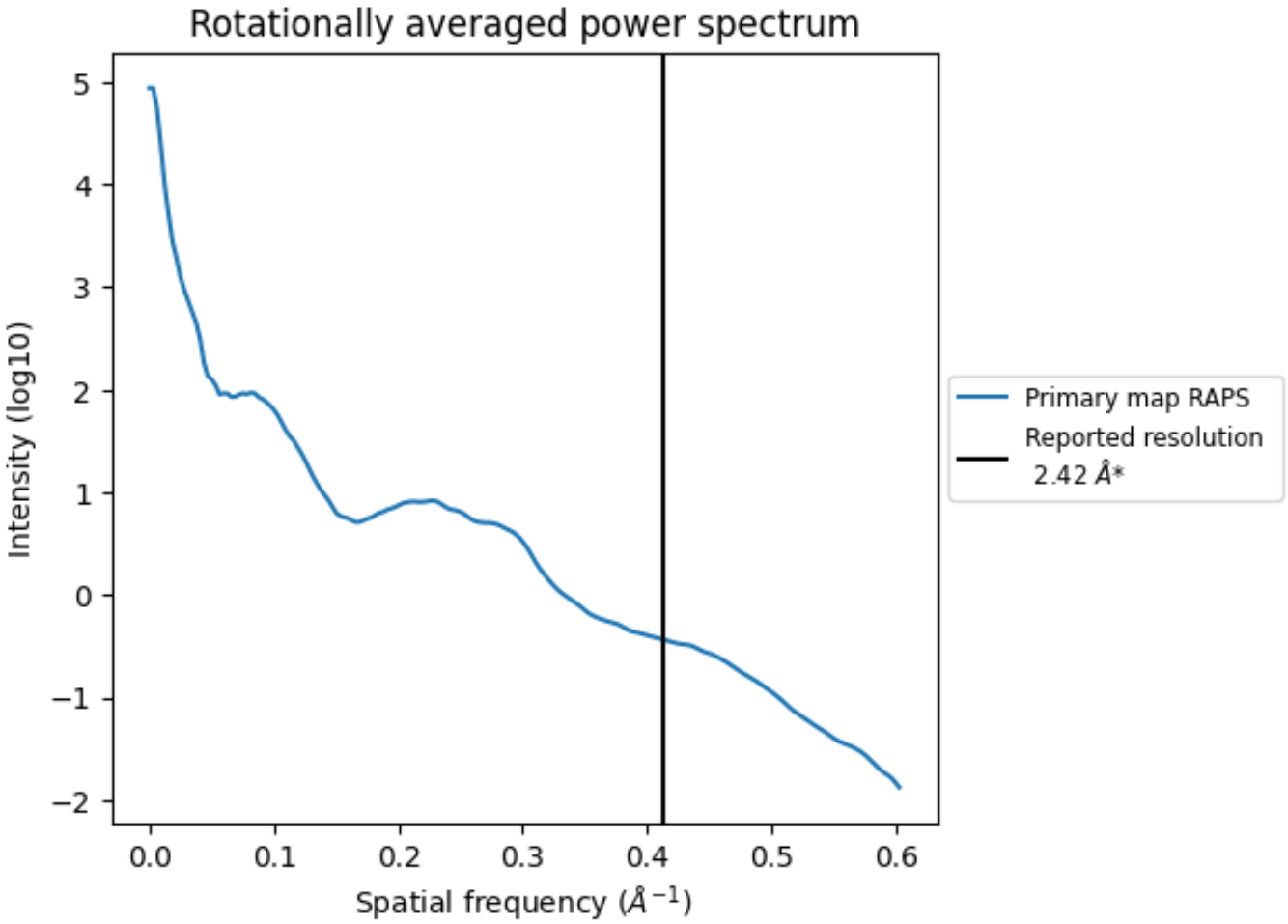
7.2 Volume estimate



The volume at the recommended contour level is 120 nm^3; this corresponds to an approximate mass of 109 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)

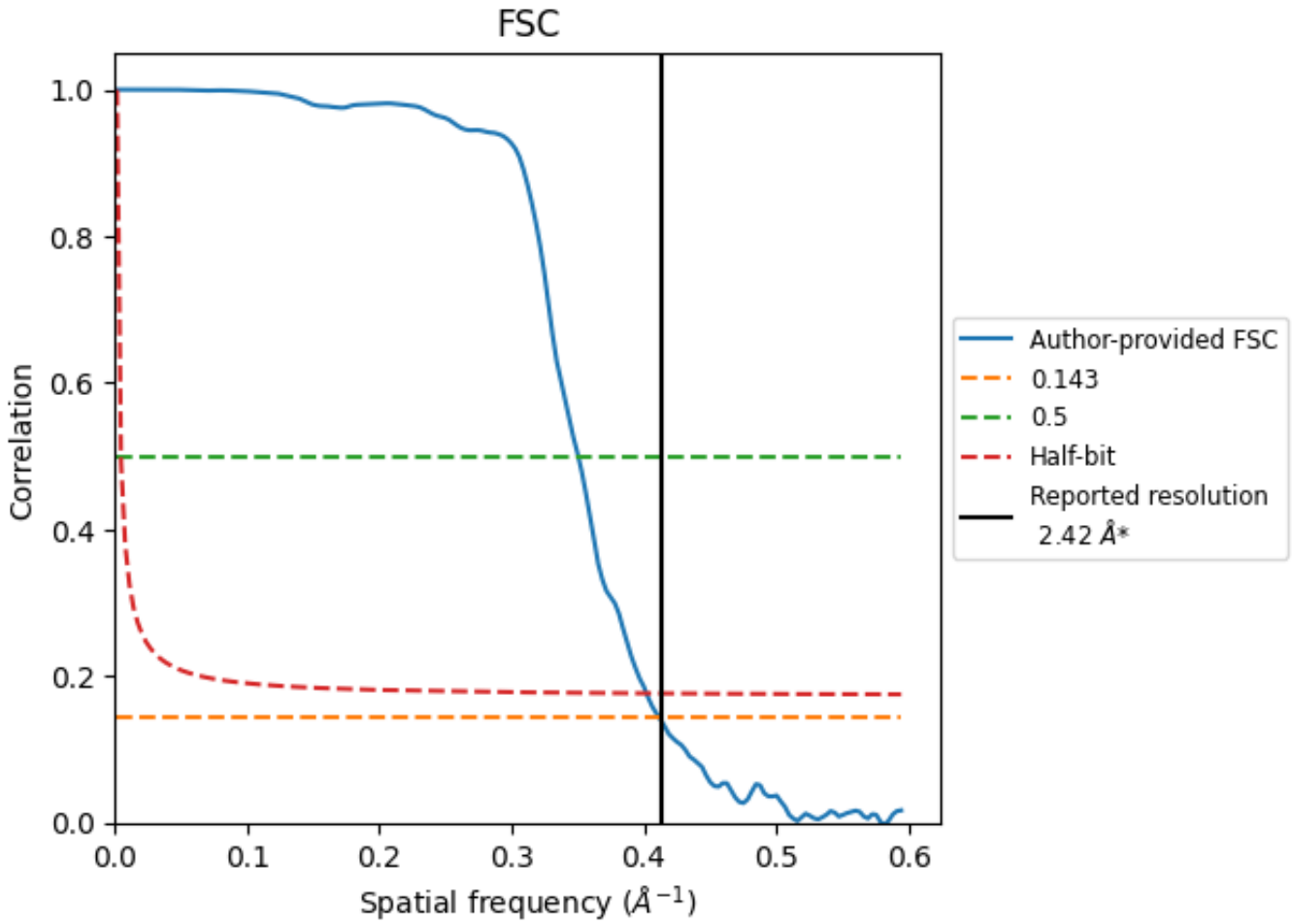


*Reported resolution corresponds to spatial frequency of 0.413 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.413 Å^{-1}

8.2 Resolution estimates (i)

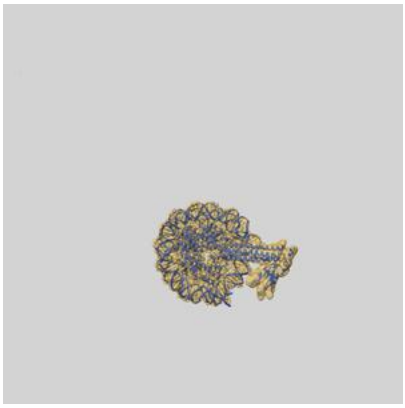
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.42	-	-
Author-provided FSC curve	2.43	2.85	2.49
Unmasked-calculated*	-	-	-

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

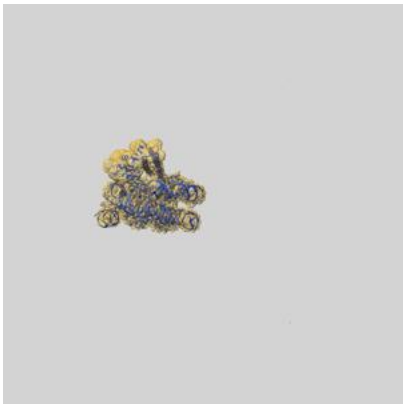
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-19513 and PDB model 8RUP. Per-residue inclusion information can be found in section 3 on page 7.

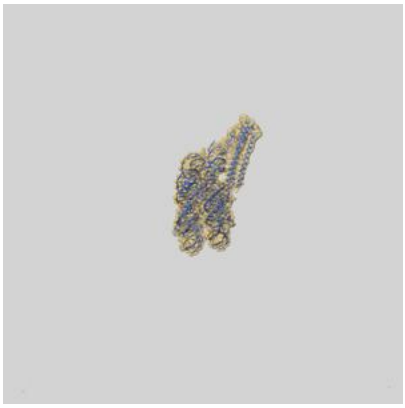
9.1 Map-model overlay



X



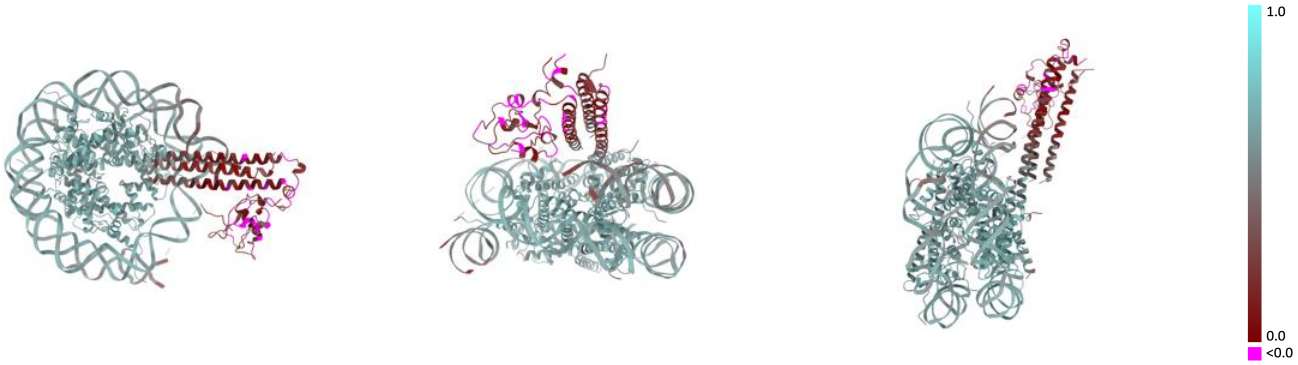
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.02 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.

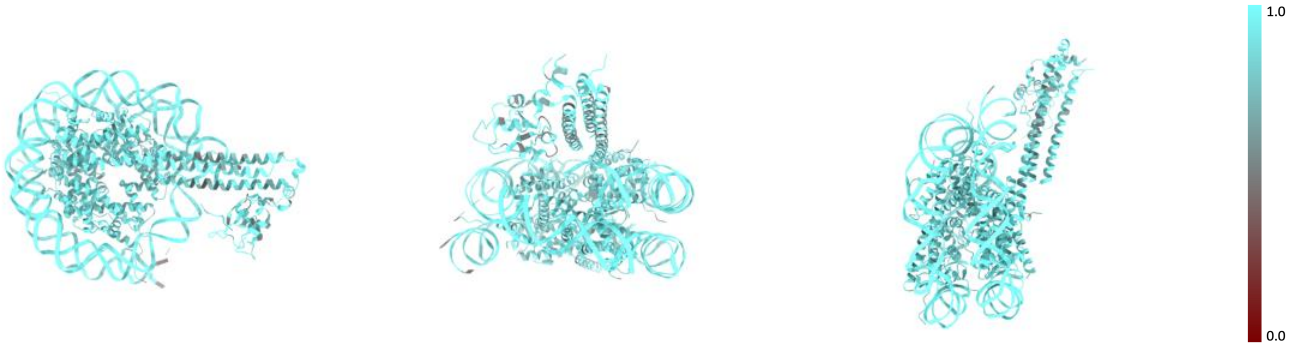
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

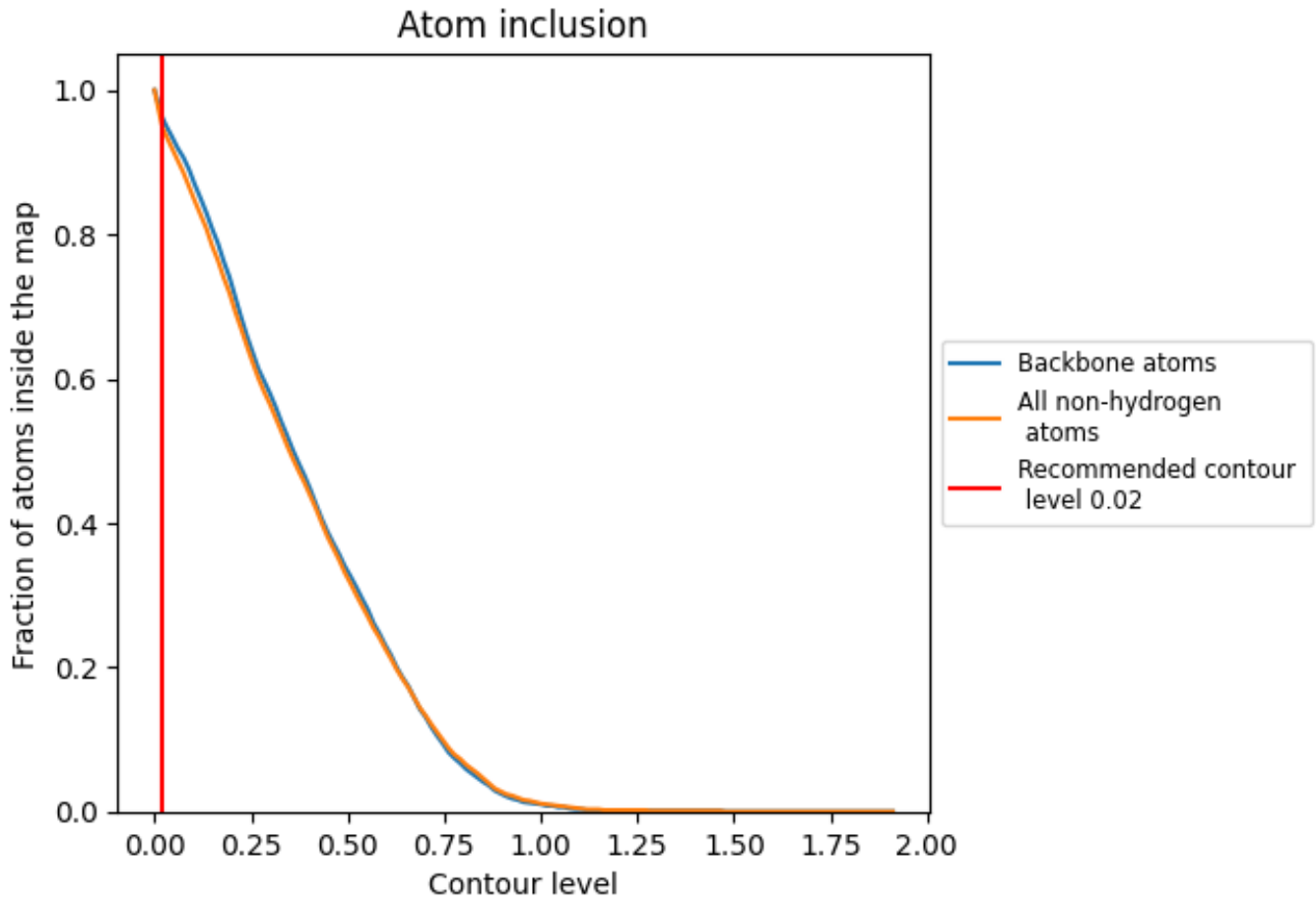
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.02).

9.4 Atom inclusion (i)



At the recommended contour level, 96% of all backbone atoms, 95% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9480	0.5510
A	0.9660	0.6410
B	0.9750	0.6560
C	0.9720	0.6490
D	0.9610	0.6380
E	0.9710	0.6290
F	0.9750	0.6680
G	0.9720	0.6440
H	0.9790	0.6480
I	0.9730	0.5720
J	0.9750	0.5750
K	0.8650	0.1400
L	0.8460	0.2650
M	0.8770	0.2570