Full wwPDB EM Validation Report

May 21, 2022 – 07:19 am BST

PDB ID	:	7PKN
EMDB ID	:	EMD-13473
Title	:	Structure of the human CCAN deltaCT complex
Authors	:	Muir, K.W.; Yatskevich, S.; Bellini, D.; Barford, D.
Deposited on	:	2021-08-25
Resolution	:	3.20 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

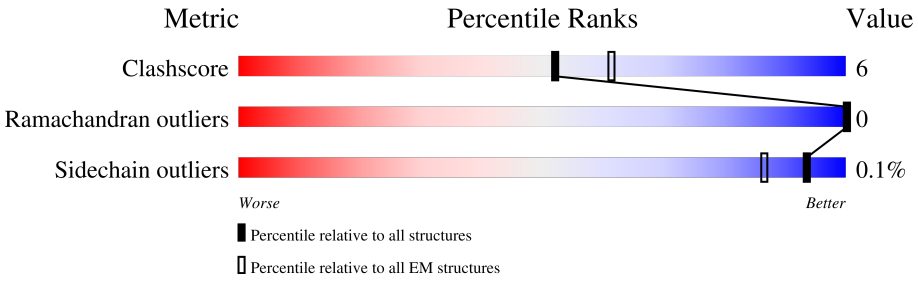
EMDB validation analysis	:	0.0.1.dev8
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

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 3.20 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	H	247	46% 11% 44%
2	I	756	44% 6% 50%
3	K	269	41% 7% 51%
4	L	344	80% 7% 13%
5	M	180	92%
6	N	339	83% 11% 6%
7	O	300	8% 53% 17% 30%
8	P	288	18% 66% 12% 22%

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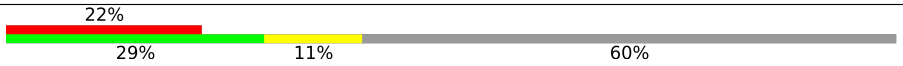
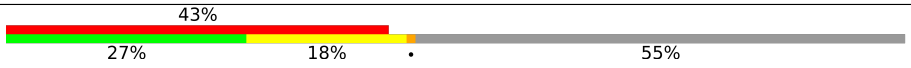
Mol	Chain	Length	Quality of chain
9	Q	268	
10	U	418	
11	R	177	

2 Entry composition (i)

There are 12 unique types of molecules in this entry. The entry contains 18616 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 Centromere protein H.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	H	139	Total	C	N	O	S	0	0
			1122	700	196	219	7		

- Molecule 2 is a protein called Centromere protein I.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	I	379	Total	C	N	O	S	0	0
			3109	2049	491	551	18		

- Molecule 3 is a protein called Centromere protein K.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	K	131	Total	C	N	O	S	0	0
			1064	661	175	220	8		

- Molecule 4 is a protein called Centromere protein L.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	L	300	Total	C	N	O	S	0	0
			2412	1571	395	432	14		

- Molecule 5 is a protein called Centromere protein M.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	172	Total	C	N	O	S	0	0
			1325	839	236	243	7		

- Molecule 6 is a protein called Centromere protein N.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	N	318	Total	C	N	O	S	0	0
			2613	1678	453	472	10		

- Molecule 7 is a protein called Centromere protein O.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	O	210	Total	C	N	O	S	0	0
			1642	1060	277	298	7		

- Molecule 8 is a protein called Centromere protein P.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	P	224	Total	C	N	O	S	0	0
			1788	1141	310	329	8		

- Molecule 9 is a protein called Centromere protein Q.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	Q	191	Total	C	N	O	S	0	0
			1526	953	258	304	11		

- Molecule 10 is a protein called Centromere protein U.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	U	166	Total	C	N	O	S	0	0
			1365	861	242	257	5		

- Molecule 11 is a protein called Centromere protein R.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	R	80	Total	C	N	O	S	0	0
			649	412	105	125	7		

- Molecule 12 is water.

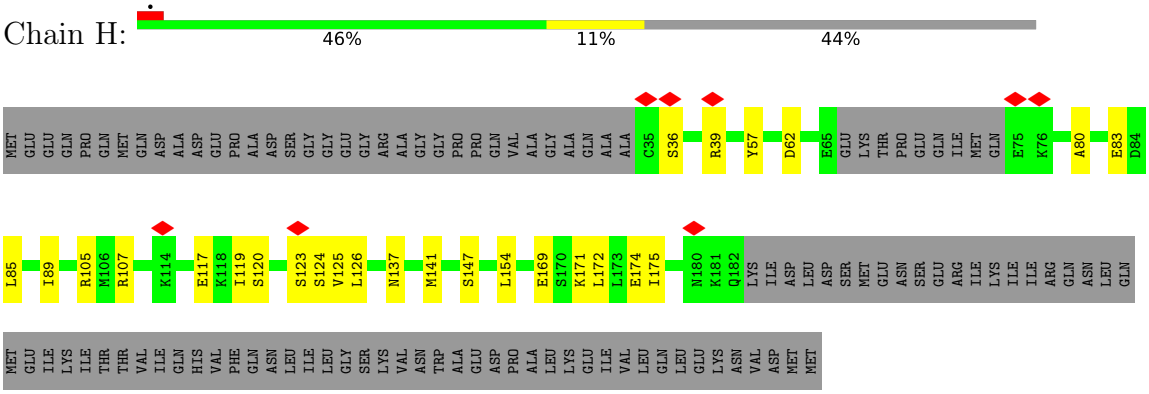
Mol	Chain	Residues	Atoms		AltConf
12	I	1	Total	O	0
			1	1	

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.

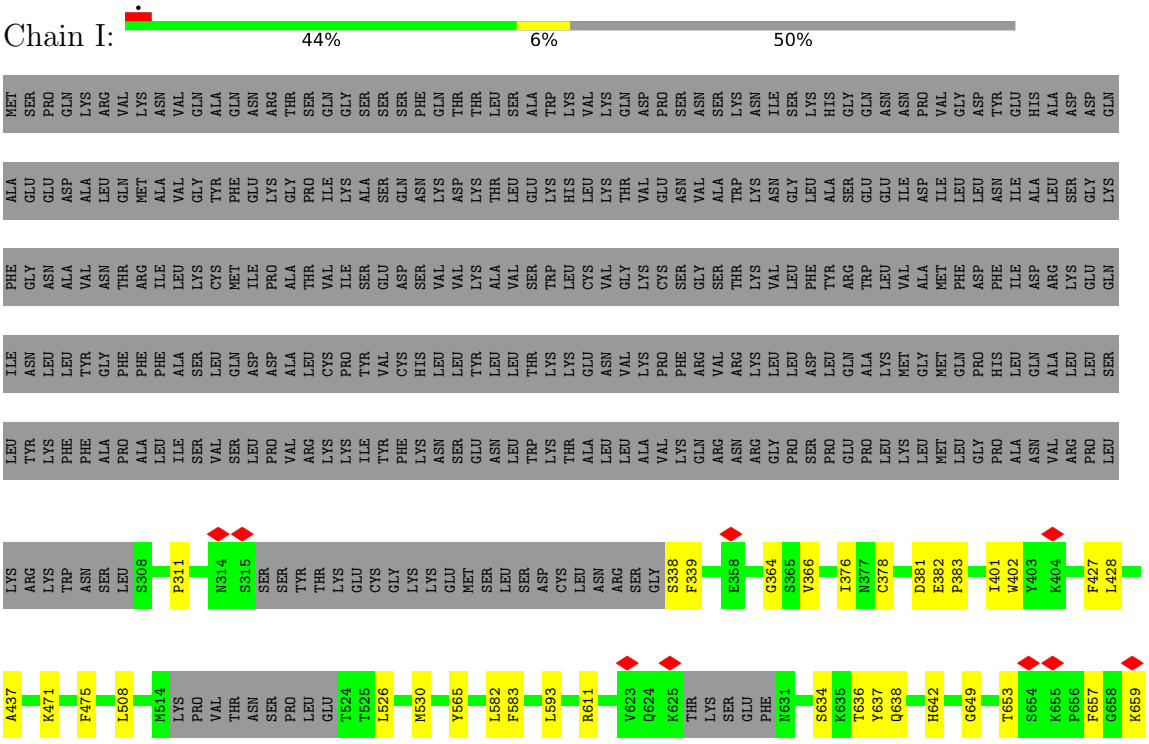
• Molecule 1: Centromere protein H

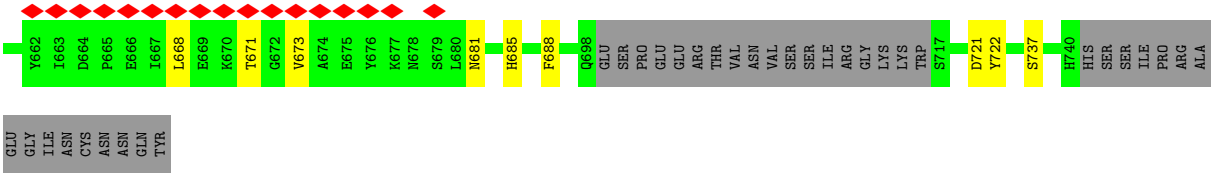
Chain H:



• Molecule 2: Centromere protein I

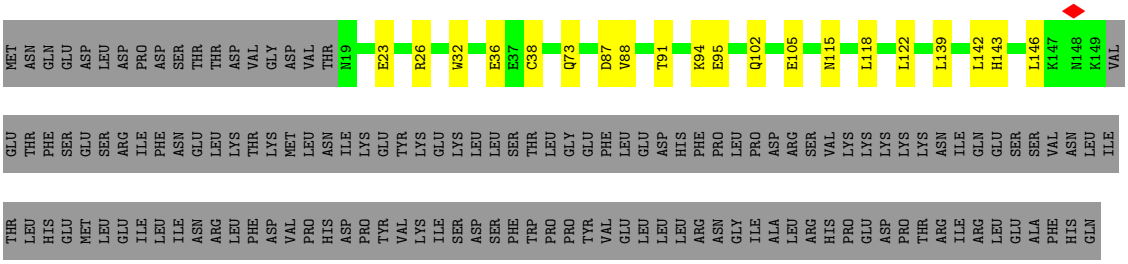
Chain I:





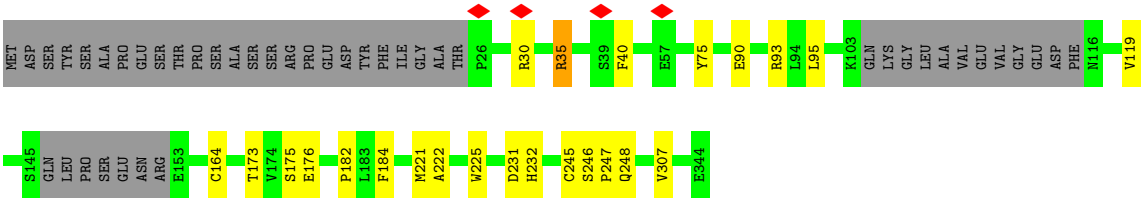
• Molecule 3: Centromere protein K

Chain K:



• Molecule 4: Centromere protein L

Chain L:

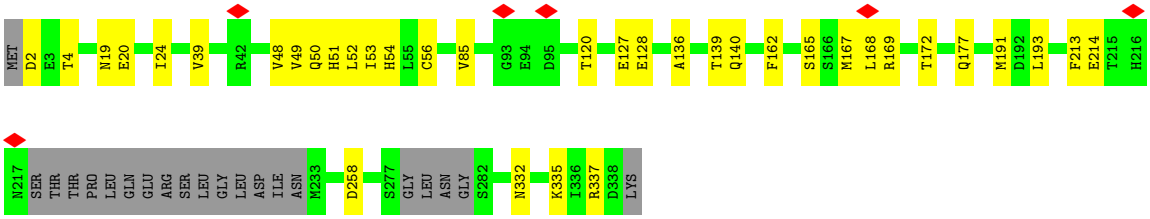


• Molecule 5: Centromere protein M

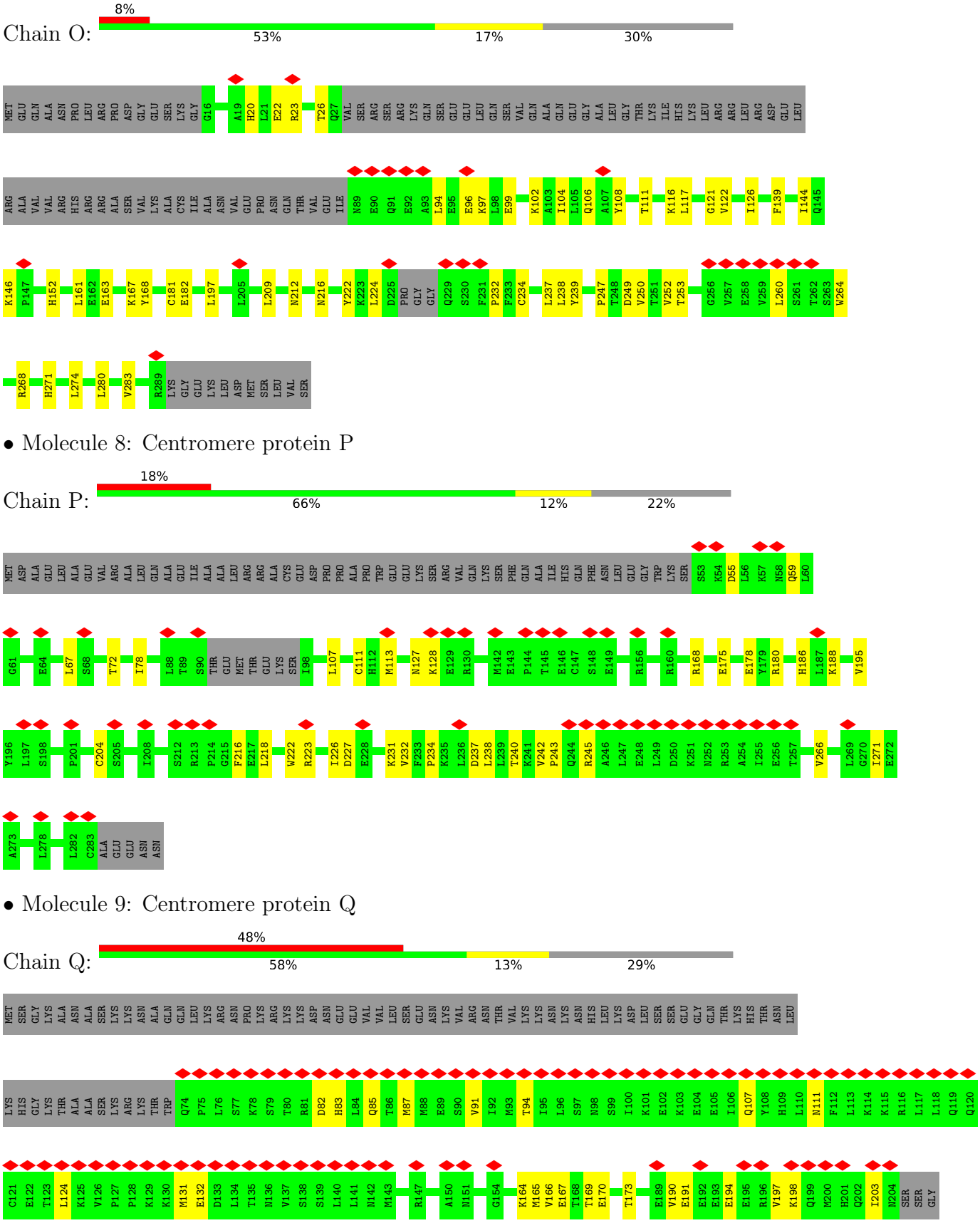
Chain M:

• Molecule 6: Centromere protein N

Chain N:

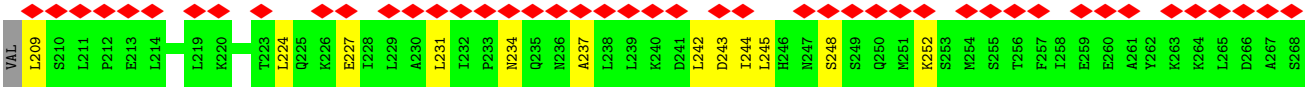


• Molecule 7: Centromere protein O

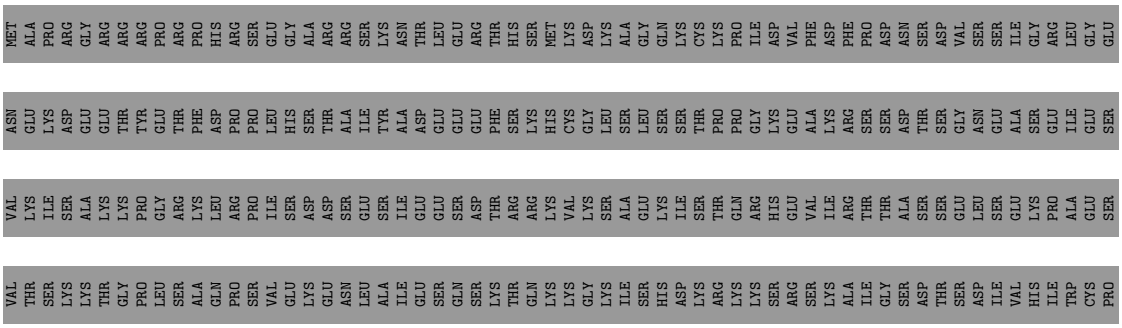


• Molecule 8: Centromere protein P

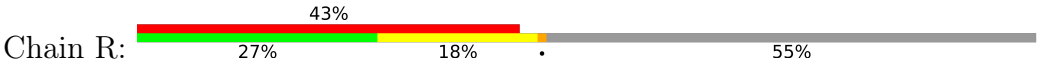
• Molecule 9: Centromere protein Q

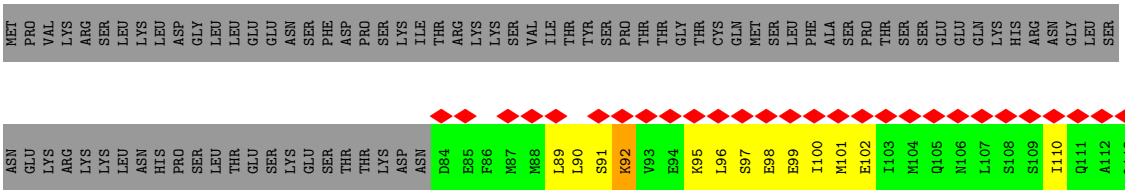


• Molecule 10: Centromere protein U



• Molecule 11: Centromere protein R





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	569755	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.057	Depositor
Minimum map value	-0.022	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.006	Depositor
Map size (Å)	292.4, 292.4, 292.4	wwPDB
Map dimensions	340, 340, 340	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.85999995, 0.85999995, 0.85999995	Depositor

5 Model quality

5.1 Standard geometry

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	H	0.25	0/1125	0.38	0/1498
2	I	0.25	0/3195	0.39	0/4336
3	K	0.25	0/1073	0.39	0/1442
4	L	0.25	0/2477	0.39	0/3360
5	M	0.25	0/1347	0.43	0/1827
6	N	0.25	0/2670	0.42	0/3606
7	O	0.25	0/1678	0.44	0/2280
8	P	0.25	0/1820	0.46	0/2451
9	Q	0.23	0/1538	0.43	0/2062
10	U	0.27	0/1383	0.44	0/1856
11	R	0.24	0/653	0.46	0/865
All	All	0.25	0/18959	0.42	0/25583

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	H	1122	0	1170	21	0
2	I	3109	0	3076	26	0
3	K	1064	0	1062	16	0
4	L	2412	0	2405	13	0
5	M	1325	0	1370	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	N	2613	0	2627	25	0
7	O	1642	0	1616	34	0
8	P	1788	0	1791	23	0
9	Q	1526	0	1586	22	0
10	U	1365	0	1396	36	0
11	R	649	0	673	24	0
12	I	1	0	0	0	0
All	All	18616	0	18772	213	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 (213) 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)
11:R:143:MET:SD	11:R:147:ASN:ND2	2.60	0.74
8:P:55:ASP:O	8:P:59:GLN:NE2	2.26	0.69
7:O:102:LYS:O	7:O:106:GLN:NE2	2.26	0.68
1:H:119:ILE:HD12	1:H:123:SER:HB3	1.76	0.68
6:N:50:GLN:O	6:N:54:HIS:ND1	2.25	0.68
7:O:167:LYS:HG3	7:O:168:TYR:HD2	1.57	0.68
7:O:234:CYS:SG	7:O:253:THR:OG1	2.52	0.68
8:P:245:ARG:HH12	10:U:375:GLU:HB2	1.58	0.68
7:O:212:ASN:ND2	7:O:216:ASN:OD1	2.28	0.67
1:H:85:LEU:O	1:H:89:ILE:HG13	1.96	0.66
1:H:117:GLU:O	1:H:120:SER:OG	2.11	0.66
1:H:175:ILE:HD12	3:K:146:LEU:HD21	1.78	0.65
8:P:226:ILE:HG12	8:P:232:VAL:HG22	1.77	0.65
2:I:649:GLY:HA2	2:I:653:THR:HB	1.78	0.65
7:O:274:LEU:HB3	7:O:283:VAL:HG11	1.79	0.65
11:R:118:GLU:HG3	11:R:119:LEU:HD12	1.79	0.64
7:O:182:GLU:OE1	8:P:168:ARG:NE	2.31	0.64
5:M:134:GLU:N	5:M:134:GLU:OE2	2.29	0.64
6:N:127:GLU:HG3	6:N:128:GLU:H	1.62	0.64
9:Q:197:VAL:HG13	9:Q:198:LYS:HG3	1.79	0.63
10:U:261:THR:HA	10:U:264:GLU:HG2	1.80	0.63
11:R:101:MET:HG2	11:R:133:LYS:HE3	1.80	0.63
8:P:237:ASP:OD1	8:P:238:LEU:N	2.31	0.63
7:O:167:LYS:HG3	7:O:168:TYR:CD2	2.35	0.62
9:Q:131:MET:HG2	9:Q:132:GLU:H	1.62	0.62
8:P:111:CYS:O	8:P:113:MET:N	2.26	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:U:384:SER:HA	10:U:387:LEU:HD13	1.83	0.61
11:R:90:LEU:HD21	11:R:140:LYS:HD2	1.83	0.61
9:Q:83:HIS:O	9:Q:87:MET:HG2	2.00	0.61
10:U:408:ASN:O	10:U:411:LEU:HB2	2.01	0.60
9:Q:198:LYS:HD3	9:Q:203:ILE:HA	1.83	0.60
6:N:120:THR:HB	6:N:136:ALA:HB3	1.84	0.60
2:I:611:ARG:NH2	2:I:737:SER:OG	2.35	0.60
8:P:180:ARG:NH1	8:P:204:CYS:SG	2.75	0.59
1:H:126:LEU:HD11	3:K:94:LYS:HG2	1.85	0.59
1:H:36:SER:HA	1:H:39:ARG:HE	1.67	0.58
11:R:97:SER:HA	11:R:100:ILE:HG12	1.86	0.57
10:U:414:LEU:HB3	11:R:151:LEU:HD23	1.86	0.57
3:K:142:LEU:HD23	3:K:146:LEU:HD23	1.85	0.57
8:P:127:ASN:OD1	8:P:128:LYS:NZ	2.37	0.57
10:U:412:GLU:HA	10:U:415:LEU:HG	1.86	0.57
2:I:366:VAL:HG11	2:I:376:ILE:HD12	1.87	0.56
1:H:62:ASP:OD2	3:K:73:GLN:NE2	2.33	0.56
2:I:381:ASP:OD2	2:I:383:PRO:HD2	2.06	0.56
4:L:222:ALA:HB2	4:L:307:VAL:HG11	1.88	0.56
1:H:137:ASN:O	1:H:141:MET:HG3	2.06	0.56
10:U:258:PHE:O	10:U:261:THR:OG1	2.24	0.56
1:H:107:ARG:HD3	2:I:671:THR:HG21	1.89	0.55
10:U:278:ILE:HA	10:U:281:PHE:CE1	2.42	0.55
10:U:402:SER:HA	10:U:405:ARG:HE	1.71	0.55
11:R:98:GLU:O	11:R:102:GLU:HG2	2.07	0.55
4:L:245:CYS:O	4:L:248:GLN:N	2.36	0.55
1:H:169:GLU:HA	2:I:378:CYS:HB2	1.88	0.54
2:I:338:SER:OG	2:I:339:PHE:N	2.39	0.54
1:H:147:SER:OG	3:K:115:ASN:OD1	2.23	0.54
9:Q:124:LEU:HD11	10:U:257:GLU:HG2	1.90	0.54
3:K:87:ASP:OD1	3:K:88:VAL:N	2.41	0.54
4:L:164:CYS:HB2	4:L:182:PRO:HD2	1.89	0.54
2:I:382:GLU:HB3	2:I:383:PRO:HD3	1.89	0.54
11:R:120:GLU:HB3	11:R:126:SER:HA	1.90	0.53
9:Q:107:GLN:O	9:Q:111:ASN:ND2	2.41	0.53
9:Q:234:ASN:HB2	9:Q:237:ALA:HB3	1.90	0.53
10:U:401:GLU:HA	10:U:404:LEU:HD12	1.91	0.53
6:N:335:LYS:HD3	6:N:337:ARG:HH22	1.73	0.53
10:U:257:GLU:O	10:U:261:THR:HG23	2.09	0.52
1:H:57:TYR:HE1	5:M:7:LEU:HD22	1.75	0.52
2:I:471:LYS:HA	2:I:475:PHE:HD2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:R:146:VAL:HA	11:R:151:LEU:HD13	1.90	0.52
6:N:177:GLN:NE2	10:U:321:ILE:O	2.35	0.51
2:I:668:LEU:HB3	2:I:673:VAL:HB	1.91	0.51
9:Q:166:VAL:HA	9:Q:169:THR:HG22	1.93	0.51
2:I:721:ASP:OD1	2:I:722:TYR:N	2.44	0.51
6:N:49:VAL:O	6:N:53:ILE:HG12	2.11	0.51
7:O:121:GLY:HA2	7:O:144:ILE:HG12	1.91	0.51
8:P:245:ARG:NH2	10:U:375:GLU:OE1	2.35	0.50
3:K:23:GLU:OE1	3:K:26:ARG:NH2	2.44	0.50
2:I:582:LEU:HD13	2:I:636:THR:HG23	1.94	0.50
7:O:126:ILE:HD11	7:O:181:CYS:HA	1.94	0.50
8:P:188:LYS:HA	8:P:195:VAL:HG11	1.93	0.50
8:P:72:THR:HG21	11:R:175:ILE:HG21	1.94	0.49
9:Q:91:VAL:HA	9:Q:94:THR:HG22	1.93	0.49
8:P:223:ARG:HH12	9:Q:224:LEU:HD13	1.75	0.49
7:O:222:TYR:OH	7:O:271:HIS:ND1	2.46	0.49
7:O:238:LEU:HB2	7:O:249:ASP:OD1	2.12	0.49
11:R:125:ILE:HG23	11:R:126:SER:H	1.78	0.49
2:I:638:GLN:O	2:I:642:HIS:ND1	2.32	0.48
10:U:301:ASN:OD1	10:U:304:ARG:NH1	2.47	0.48
10:U:265:HIS:HD2	10:U:268:ARG:HH22	1.61	0.48
7:O:108:TYR:O	7:O:111:THR:HG22	2.13	0.48
1:H:119:ILE:O	1:H:119:ILE:HG13	2.13	0.48
1:H:175:ILE:HD11	3:K:143:HIS:ND1	2.29	0.48
2:I:685:HIS:HB3	2:I:688:PHE:HD2	1.77	0.48
4:L:246:SER:OG	4:L:247:PRO:HD3	2.13	0.48
7:O:22:GLU:O	7:O:26:THR:HG23	2.14	0.48
11:R:91:SER:O	11:R:95:LYS:HG2	2.13	0.48
7:O:20:HIS:HA	7:O:23:ARG:HG2	1.95	0.47
7:O:117:LEU:HD23	7:O:122:VAL:HG22	1.96	0.47
9:Q:243:ASP:OD2	9:Q:244:ILE:N	2.47	0.47
10:U:325:ASP:O	10:U:329:ARG:HG2	2.13	0.47
11:R:149:GLN:HB3	11:R:151:LEU:HD12	1.94	0.47
7:O:252:VAL:O	7:O:268:ARG:NH2	2.47	0.47
4:L:95:LEU:HD23	4:L:119:VAL:HG11	1.97	0.47
8:P:266:VAL:HG23	8:P:271:ILE:HA	1.95	0.47
9:Q:164:LYS:HG3	9:Q:165:MET:N	2.30	0.47
10:U:404:LEU:HA	10:U:407:ILE:HG22	1.97	0.47
1:H:124:SER:OG	1:H:125:VAL:N	2.48	0.47
6:N:140:GLN:OE1	10:U:317:ARG:NH2	2.43	0.47
6:N:258:ASP:OD1	6:N:258:ASP:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:O:96:GLU:O	7:O:99:GLU:HG3	2.14	0.47
9:Q:191:GLU:HA	9:Q:194:GLU:HG3	1.96	0.47
2:I:437:ALA:HB1	4:L:35:ARG:HE	1.81	0.46
10:U:278:ILE:HD13	10:U:281:PHE:CE1	2.50	0.46
11:R:96:LEU:HD12	11:R:99:GLU:OE2	2.15	0.46
3:K:32:TRP:HZ3	5:M:10:LEU:HD13	1.80	0.46
6:N:85:VAL:HG23	6:N:191:MET:HB2	1.98	0.46
9:Q:197:VAL:HG23	10:U:355:ALA:HB2	1.97	0.46
10:U:262:HIS:O	10:U:266:GLN:HG2	2.16	0.46
7:O:146:LYS:HE3	11:R:169:TYR:HD2	1.80	0.46
4:L:175:SER:OG	4:L:176:GLU:N	2.49	0.46
9:Q:190:VAL:HG23	10:U:348:LYS:HG3	1.97	0.46
11:R:97:SER:HB2	11:R:101:MET:HE1	1.96	0.46
8:P:78:ILE:HD11	8:P:107:LEU:HD22	1.98	0.46
8:P:227:ASP:OD1	8:P:231:LYS:N	2.49	0.45
9:Q:167:GLU:O	9:Q:170:GLU:HG3	2.16	0.45
9:Q:227:GLU:O	9:Q:231:LEU:HB2	2.15	0.45
11:R:96:LEU:HA	11:R:99:GLU:OE2	2.16	0.45
8:P:111:CYS:C	8:P:113:MET:H	2.12	0.45
10:U:396:THR:O	10:U:400:ALA:N	2.50	0.45
11:R:89:LEU:O	11:R:92:LYS:HG3	2.16	0.45
9:Q:82:ASP:O	9:Q:85:GLN:HG3	2.17	0.45
11:R:110:ILE:O	11:R:114:GLU:HG3	2.17	0.45
11:R:96:LEU:O	11:R:100:ILE:HG23	2.17	0.44
7:O:163:GLU:O	7:O:167:LYS:HG2	2.17	0.44
9:Q:170:GLU:HA	9:Q:173:THR:HG22	2.00	0.44
10:U:413:LYS:HD3	10:U:413:LYS:HA	1.86	0.44
2:I:401:ILE:HB	2:I:402:TRP:CE3	2.53	0.44
6:N:39:VAL:HG22	6:N:51:HIS:ND1	2.32	0.44
10:U:313:ILE:O	10:U:317:ARG:HG3	2.17	0.44
11:R:138:LYS:HB3	11:R:138:LYS:HE2	1.75	0.44
2:I:311:PRO:HD2	2:I:364:GLY:HA3	2.00	0.44
2:I:634:SER:HA	2:I:637:TYR:HD1	1.82	0.44
2:I:657:PHE:HZ	2:I:668:LEU:HD11	1.82	0.44
7:O:94:LEU:HA	7:O:97:LYS:HE2	2.00	0.44
8:P:175:GLU:O	8:P:178:GLU:HG2	2.18	0.43
6:N:2:ASP:OD2	6:N:4:THR:OG1	2.24	0.43
7:O:197:LEU:HD22	7:O:280:LEU:HD21	1.99	0.43
1:H:57:TYR:HB3	3:K:38:CYS:SG	2.58	0.43
4:L:30:ARG:HG2	4:L:173:THR:HG22	2.00	0.43
7:O:237:LEU:HG	7:O:250:VAL:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:U:326:GLU:O	10:U:330:LEU:HD23	2.18	0.43
10:U:407:ILE:O	10:U:411:LEU:N	2.49	0.43
2:I:565:TYR:CE1	2:I:583:PHE:HB3	2.53	0.43
8:P:216:PHE:HE2	8:P:242:VAL:HG23	1.83	0.43
1:H:105:ARG:HG3	2:I:593:LEU:HB3	2.00	0.43
6:N:168:LEU:O	6:N:172:THR:N	2.51	0.43
8:P:216:PHE:CZ	8:P:243:PRO:HD2	2.53	0.43
10:U:322:GLU:O	10:U:326:GLU:OE1	2.37	0.43
4:L:40:PHE:O	5:M:115:HIS:ND1	2.37	0.43
6:N:19:ASN:OD1	6:N:19:ASN:N	2.51	0.43
7:O:104:ILE:HG21	8:P:67:LEU:HD11	2.00	0.43
8:P:186:HIS:CE1	9:Q:209:LEU:HD21	2.54	0.43
10:U:358:LEU:HD12	10:U:358:LEU:HA	1.92	0.43
5:M:106:ARG:HG3	5:M:106:ARG:HH11	1.84	0.43
7:O:161:LEU:H	7:O:161:LEU:HD23	1.83	0.43
2:I:508:LEU:HD23	2:I:508:LEU:HA	1.86	0.42
7:O:146:LYS:HE3	11:R:169:TYR:CD2	2.54	0.42
7:O:232:PRO:HD2	7:O:264:TRP:CH2	2.54	0.42
11:R:148:LYS:HA	11:R:148:LYS:HD3	1.79	0.42
6:N:48:VAL:O	6:N:52:LEU:HD23	2.19	0.42
6:N:165:SER:O	6:N:165:SER:OG	2.34	0.42
10:U:269:ILE:HG23	10:U:275:LYS:HE3	2.02	0.42
11:R:135:GLU:O	11:R:139:THR:HG23	2.19	0.42
4:L:90:GLU:HA	4:L:93:ARG:HG2	2.01	0.42
5:M:35:MET:HG2	5:M:147:VAL:HG23	2.00	0.42
7:O:239:TYR:CD1	7:O:247:PRO:HA	2.54	0.42
1:H:172:LEU:HD13	3:K:139:LEU:HD11	2.01	0.42
3:K:36:GLU:HG2	5:M:10:LEU:HD11	2.02	0.42
6:N:20:GLU:O	6:N:24:ILE:HG12	2.19	0.42
8:P:218:LEU:HD22	8:P:240:THR:HG22	2.02	0.42
2:I:427:PHE:O	2:I:428:LEU:HB2	2.20	0.42
4:L:231:ASP:OD1	4:L:232:HIS:ND1	2.46	0.42
10:U:310:ILE:HA	10:U:313:ILE:HG12	2.02	0.42
6:N:53:ILE:HA	6:N:56:CYS:SG	2.59	0.42
6:N:85:VAL:HG12	6:N:162:PHE:CD2	2.55	0.42
3:K:91:THR:O	3:K:95:GLU:HG2	2.19	0.41
6:N:213:PHE:CD2	7:O:139:PHE:HB2	2.56	0.41
2:I:401:ILE:HB	2:I:402:TRP:HE3	1.83	0.41
4:L:221:MET:HE3	4:L:225:TRP:CH2	2.55	0.41
7:O:209:LEU:HD23	7:O:209:LEU:H	1.84	0.41
6:N:213:PHE:HE1	7:O:152:HIS:CD2	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:214:GLU:HB3	7:O:116:LYS:HD3	2.03	0.41
10:U:261:THR:HB	10:U:265:HIS:CE1	2.55	0.41
1:H:80:ALA:O	1:H:83:GLU:HG3	2.21	0.41
8:P:222:TRP:HE1	8:P:234:PRO:HB2	1.86	0.41
6:N:335:LYS:HB3	6:N:337:ARG:HH12	1.85	0.41
3:K:102:GLN:O	3:K:105:GLU:HG2	2.20	0.41
9:Q:248:SER:O	9:Q:252:LYS:HG2	2.20	0.41
1:H:171:LYS:O	1:H:174:GLU:HG2	2.21	0.41
3:K:118:LEU:HD23	3:K:118:LEU:HA	1.93	0.41
6:N:193:LEU:HD12	6:N:332:ASN:HB2	2.03	0.41
6:N:139:THR:OG1	6:N:140:GLN:N	2.54	0.41
7:O:260:LEU:HD23	7:O:260:LEU:H	1.86	0.41
10:U:411:LEU:HD23	10:U:411:LEU:HA	1.92	0.41
6:N:167:MET:HA	6:N:169:ARG:NH1	2.36	0.40
10:U:256:PRO:O	10:U:260:LYS:HG2	2.22	0.40
1:H:154:LEU:HD21	3:K:122:LEU:HA	2.03	0.40
9:Q:242:LEU:O	9:Q:245:LEU:HG	2.22	0.40
2:I:526:LEU:O	2:I:530:MET:HG2	2.21	0.40
4:L:75:TYR:HB2	4:L:184:PHE:HD2	1.87	0.40
7:O:122:VAL:HG23	7:O:144:ILE:HD11	2.04	0.40
10:U:368:TYR:O	10:U:370:ASP:N	2.54	0.40
2:I:681:ASN:O	2:I:685:HIS:HB2	2.22	0.40
7:O:224:LEU:HD23	7:O:224:LEU:H	1.86	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	H	135/247 (55%)	129 (96%)	6 (4%)	0	100	100
2	I	369/756 (49%)	357 (97%)	12 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	K	129/269 (48%)	126 (98%)	3 (2%)	0	100	100
4	L	294/344 (86%)	292 (99%)	2 (1%)	0	100	100
5	M	170/180 (94%)	169 (99%)	1 (1%)	0	100	100
6	N	312/339 (92%)	300 (96%)	12 (4%)	0	100	100
7	O	204/300 (68%)	201 (98%)	3 (2%)	0	100	100
8	P	220/288 (76%)	212 (96%)	8 (4%)	0	100	100
9	Q	187/268 (70%)	181 (97%)	6 (3%)	0	100	100
10	U	164/418 (39%)	162 (99%)	2 (1%)	0	100	100
11	R	76/177 (43%)	72 (95%)	4 (5%)	0	100	100
All	All	2260/3586 (63%)	2201 (97%)	59 (3%)	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	H	126/224 (56%)	126 (100%)	0	100	100
2	I	353/691 (51%)	352 (100%)	1 (0%)	92	96
3	K	123/260 (47%)	123 (100%)	0	100	100
4	L	268/306 (88%)	267 (100%)	1 (0%)	91	95
5	M	151/158 (96%)	151 (100%)	0	100	100
6	N	287/311 (92%)	287 (100%)	0	100	100
7	O	177/263 (67%)	177 (100%)	0	100	100
8	P	197/259 (76%)	197 (100%)	0	100	100
9	Q	179/248 (72%)	179 (100%)	0	100	100
10	U	152/379 (40%)	152 (100%)	0	100	100
11	R	75/166 (45%)	74 (99%)	1 (1%)	69	87
All	All	2088/3265 (64%)	2085 (100%)	3 (0%)	93	98

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	I	659	LYS
4	L	35	ARG
11	R	92	LYS

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

Mol	Chain	Res	Type
7	O	106	GLN
8	P	59	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides 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.

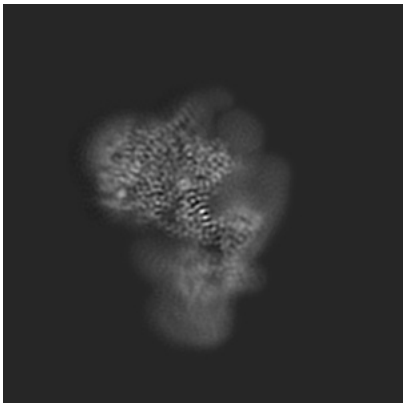
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-13473. These allow visual inspection of the internal detail of the map and identification of artifacts.

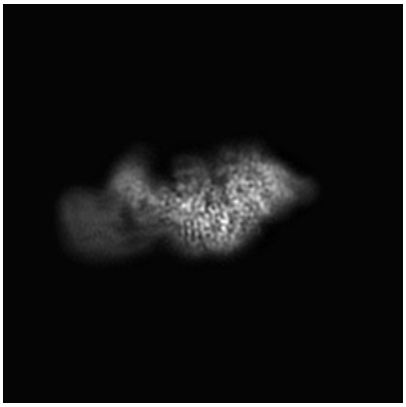
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

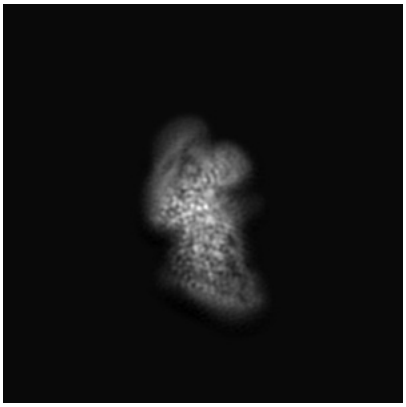
6.1.1 Primary map



X



Y

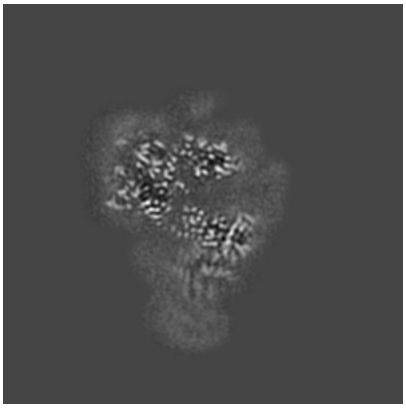


Z

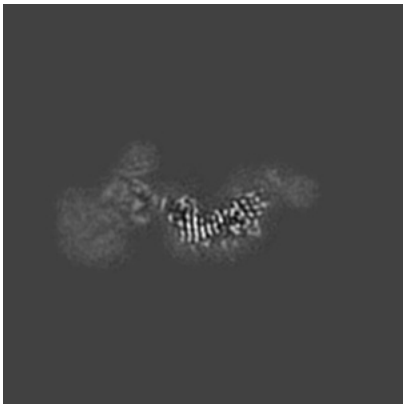
The images above show the map projected in three orthogonal directions.

6.2 Central slices

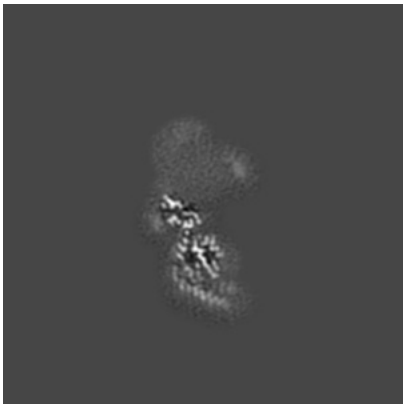
6.2.1 Primary map



X Index: 170



Y Index: 170

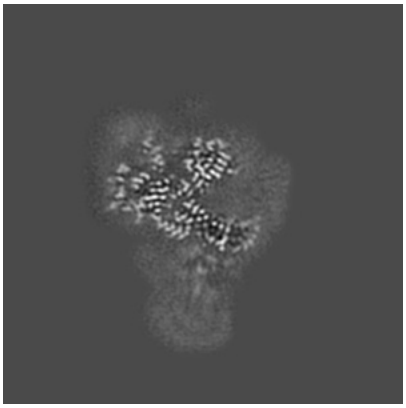


Z Index: 170

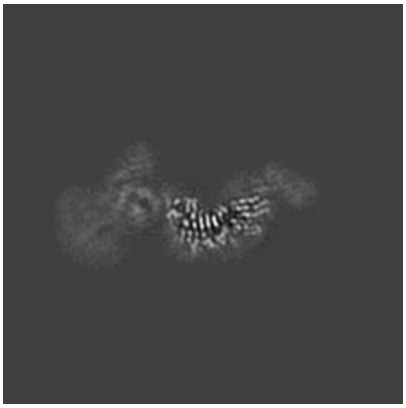
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

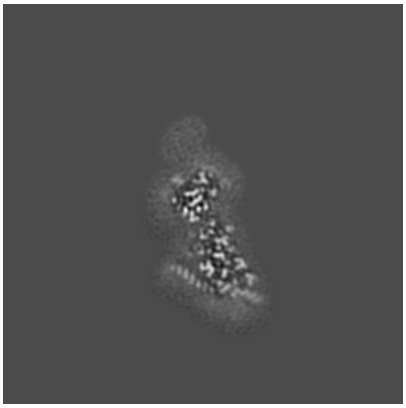
6.3.1 Primary map



X Index: 165



Y Index: 167



Z Index: 202

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

6.4 Orthogonal surface views (i)

6.4.1 Primary map

X

Y

Z

The images above show the 3D surface view of the map at the recommended contour level 0.006. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

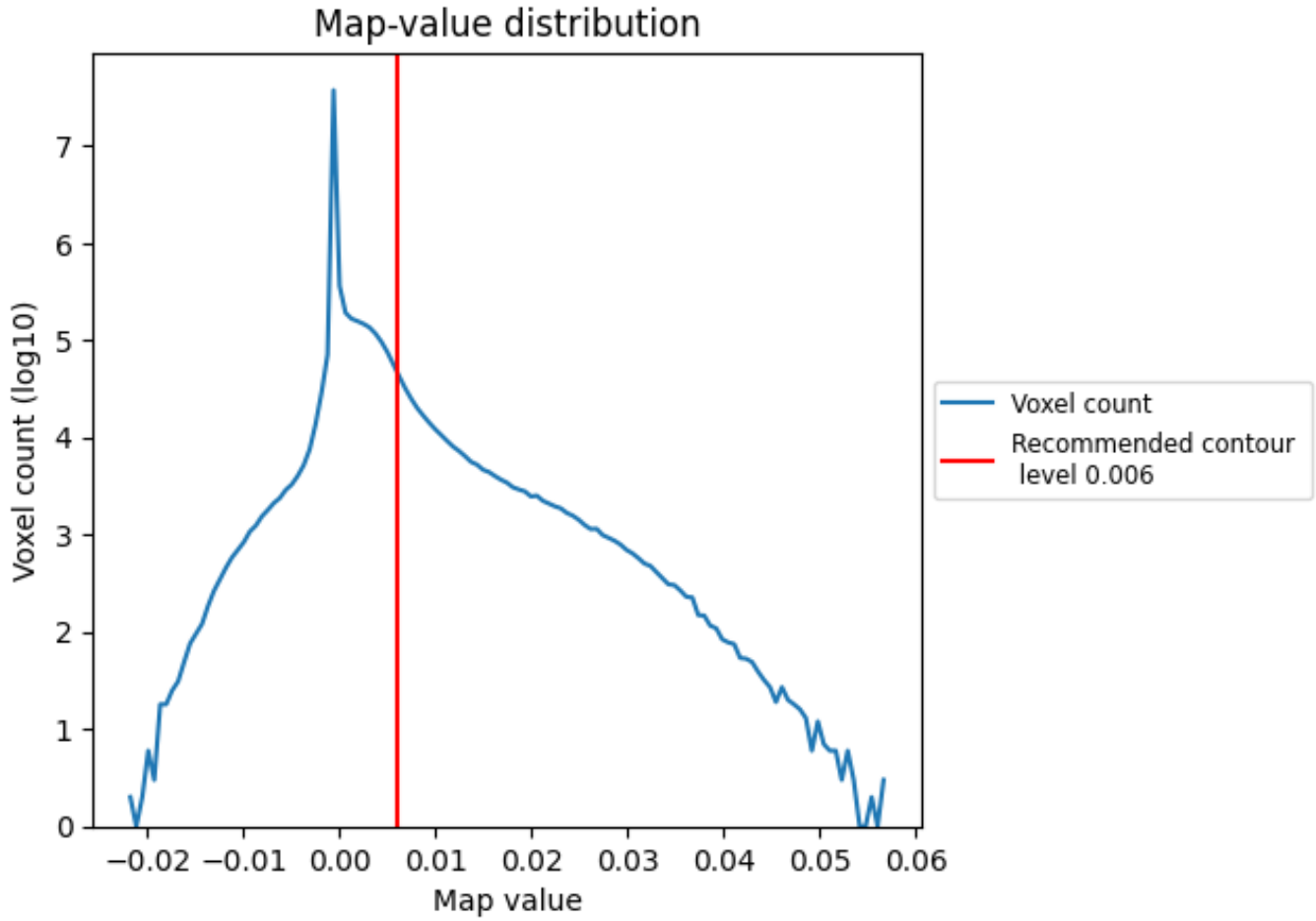
6.5 Mask visualisation (i)

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

7 Map analysis

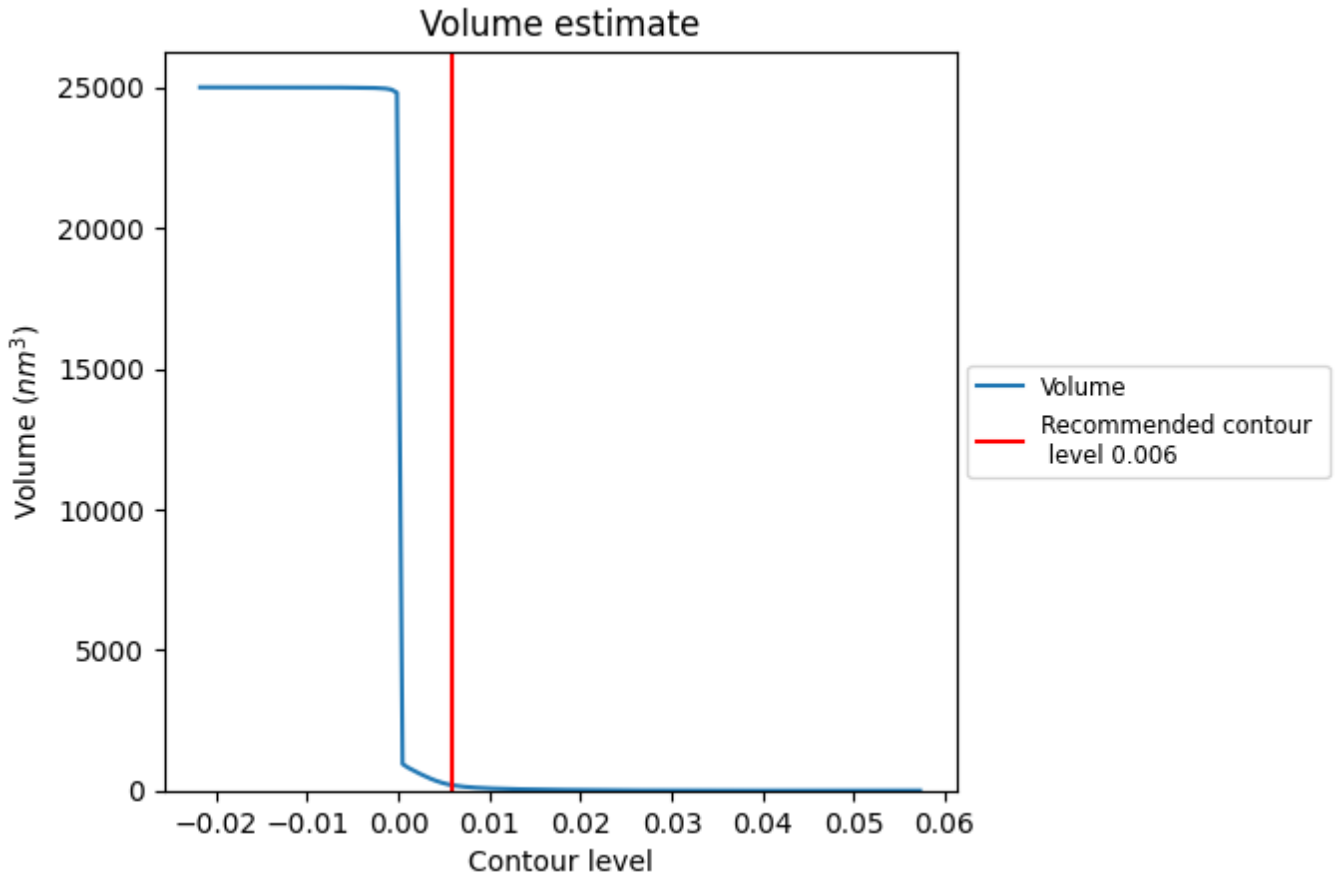
This section contains the results of statistical analysis of the map.

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.

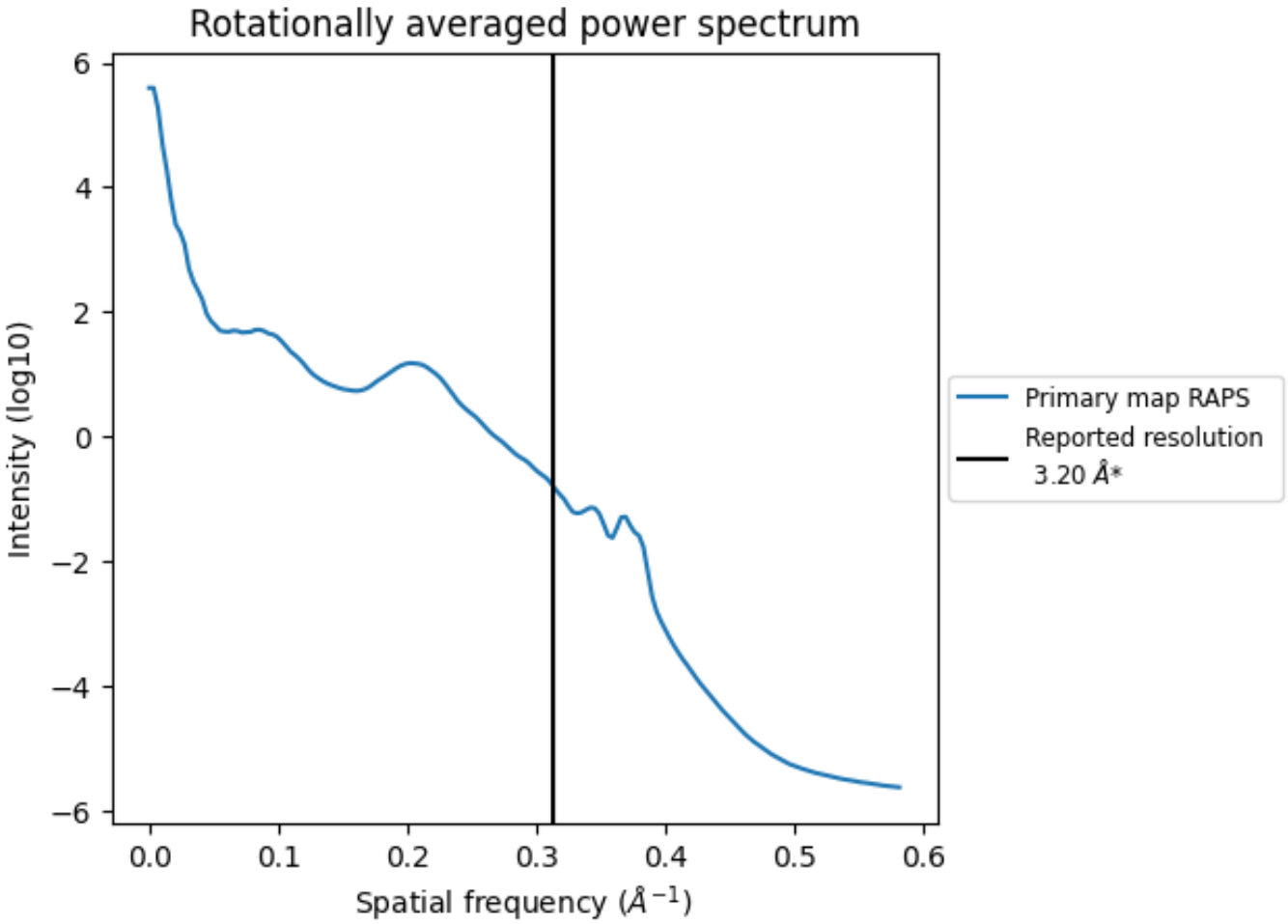
7.2 Volume estimate



The volume at the recommended contour level is 188 nm^3; this corresponds to an approximate mass of 170 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)

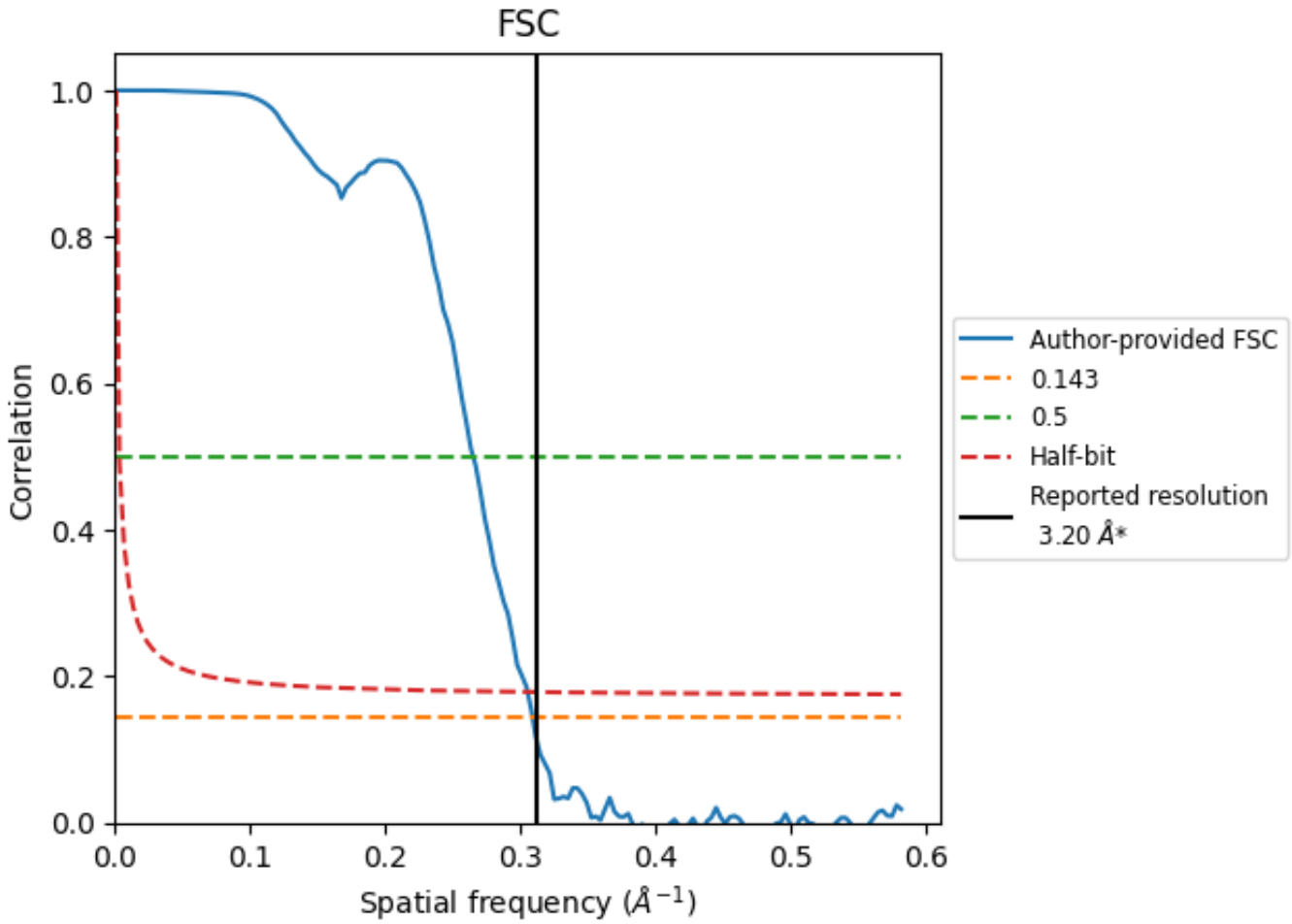


*Reported resolution corresponds to spatial frequency of 0.312 Å^{-1}

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 0.312 Å^{-1}

8.2 Resolution estimates (i)

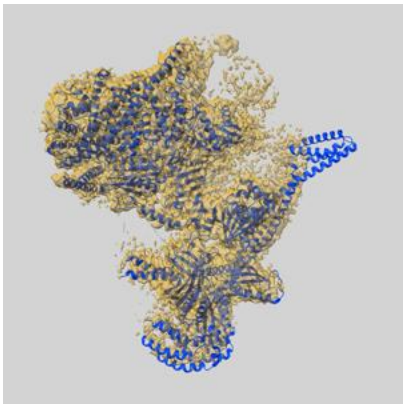
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.24	3.78	3.28
Unmasked-calculated*	-	-	-

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

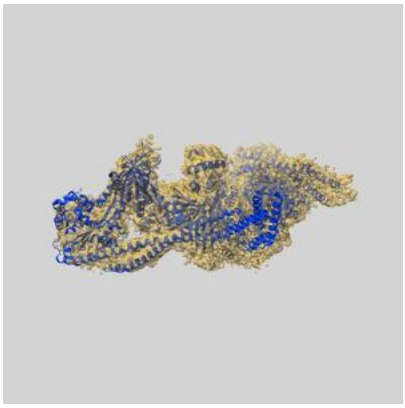
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-13473 and PDB model 7PKN. Per-residue inclusion information can be found in section 3 on page 6.

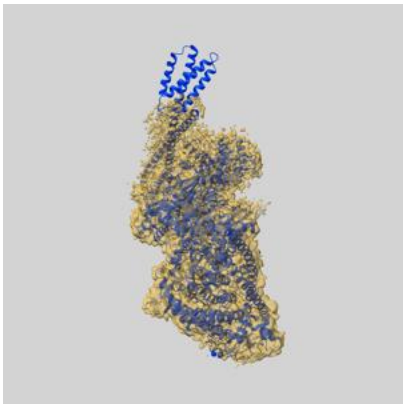
9.1 Map-model overlay



X



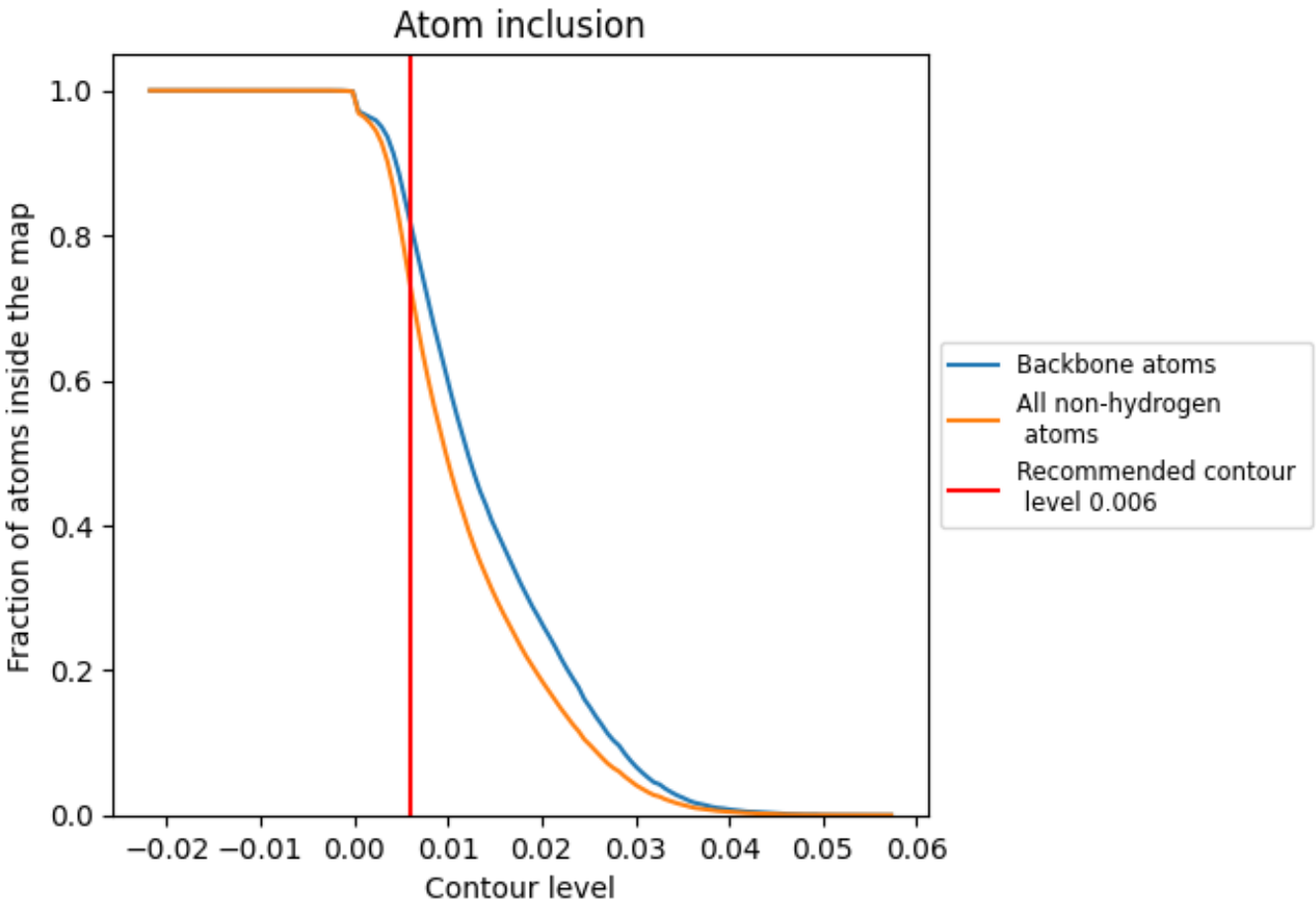
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.006 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 Atom inclusion (i)



At the recommended contour level, 82% of all backbone atoms, 73% of all non-hydrogen atoms, are inside the map.