



Full wwPDB EM Validation Report ⓘ

Nov 29, 2022 – 09:31 AM JST

PDB ID : 7W9V
EMDB ID : EMD-32373
Title : Cryo-EM structure of nucleosome in complex with p300 acetyltransferase catalytic core (complex I)
Authors : Hatazawa, S.; Liu, J.; Takizawa, Y.; Zandian, M.; Negishi, L.; Kutateladze, T.G.; Kurumizaka, H.
Deposited on : 2021-12-10
Resolution : 3.95 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:

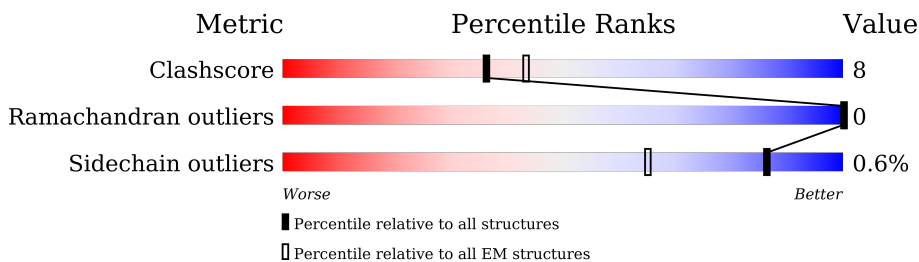
EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

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 3.95 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	135	57% 15% 28%
1	E	135	58% 14% 28%
2	B	106	60% 18% 22%
2	F	106	50% 24% 26%
3	C	133	70% 8% 23%
3	G	133	72% 7% 21%
4	D	129	62% 12% 26%
4	H	129	60% 12% 28%

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Mol	Chain	Length	Quality of chain
5	I	145	 52% 48%
6	J	145	 55% 45%
7	K	638	 59% 77% 9% 14%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 18546 atoms, of which 2136 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Histone H3.1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
1	A	97	Total	C	H	N	O	S	0	0
			1008	505	207	155	137	4		
1	E	97	Total	C	H	N	O	S	0	0
			1008	503	208	155	138	4		

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	B	83	Total	C	H	N	O	S	0	0
			841	418	179	129	114	1		
2	F	78	Total	C	H	N	O	S	0	0
			785	391	166	120	107	1		

There are 6 discrepancies between the modelled and reference sequences:

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
3	C	103	Total	C	H	N	O	0	0
			996	502	200	155	139		
3	G	105	Total	C	H	N	O	0	0
			1015	511	205	158	141		

There are 6 discrepancies between the modelled and reference sequences:

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
4	D	95	Total	C	H	N	O	S	0	0
			937	468	191	136	140	2		
4	H	93	Total	C	H	N	O	S	0	0
			906	456	181	130	137	2		

There are 6 discrepancies between the modelled and reference sequences:

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

- Molecule 5 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace	
5	I	145	Total	C	H	N	O	P	0	0
			3246	1404	294	537	867	144		

- Molecule 6 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace	
6	J	145	Total	C	H	N	O	P	0	0
			3292	1416	305	558	869	144		

- Molecule 7 is a protein called Histone acetyltransferase p300.

Mol	Chain	Residues	Atoms				AltConf	Trace	
7	K	551	Total	C	N	O	S	0	0
			4512	2884	769	826	33		


There are 14 discrepancies between the modelled and reference sequences:

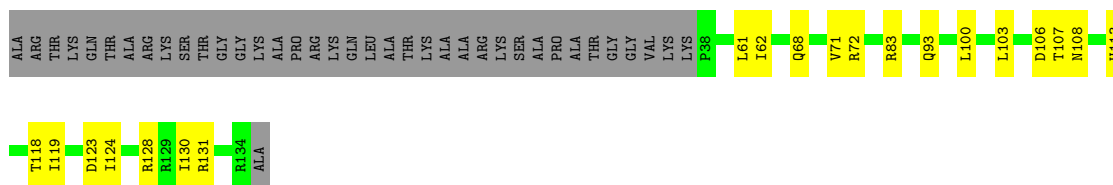
Chain	Residue	Modelled	Actual	Comment	Reference
K	1467	PHE	TYR	engineered mutation	UNP Q09472
K	1520	SER	-	linker	UNP Q09472
K	1521	GLY	-	linker	UNP Q09472
K	1522	GLY	-	linker	UNP Q09472
K	1523	SER	-	linker	UNP Q09472
K	1580	GLY	-	linker	UNP Q09472
K	1721	HIS	-	expression tag	UNP Q09472
K	1722	HIS	-	expression tag	UNP Q09472
K	1723	HIS	-	expression tag	UNP Q09472
K	1724	HIS	-	expression tag	UNP Q09472
K	1725	HIS	-	expression tag	UNP Q09472
K	1726	HIS	-	expression tag	UNP Q09472
K	1727	HIS	-	expression tag	UNP Q09472
K	1728	HIS	-	expression tag	UNP Q09472

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.

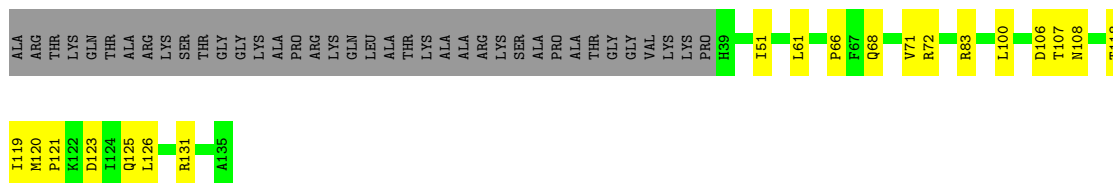
- Molecule 1: Histone H3.1

Chain A:  57% 15% 28%



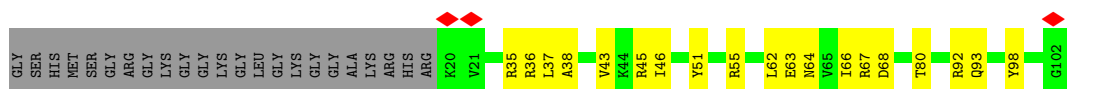
- Molecule 1: Histone H3.1

Chain E:  58% 14% 28%



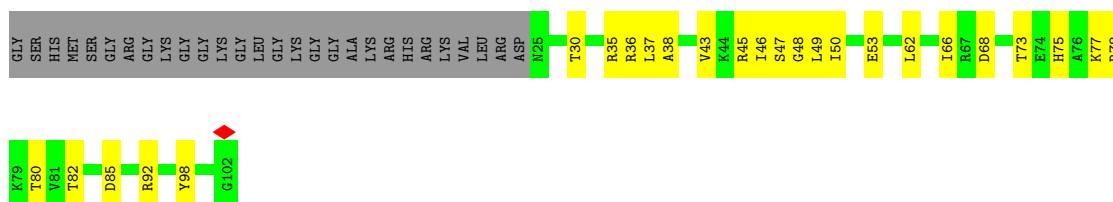
- Molecule 2: Histone H4

Chain B:  60% 18% 22%



- Molecule 2: Histone H4

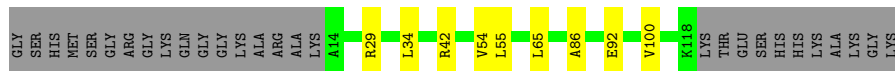
Chain F:  50% 24% 26%



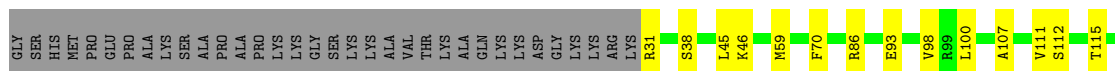
• Molecule 3: Histone H2A type 1-B/E



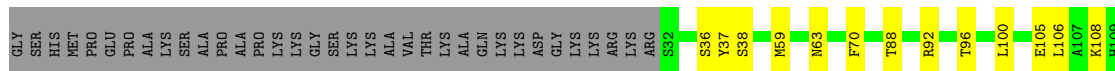
• Molecule 3: Histone H2A type 1-B/E



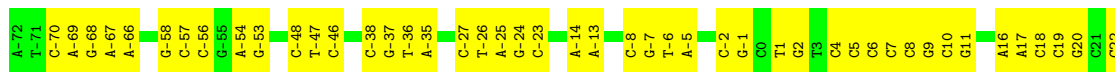
• Molecule 4: Histone H2B type 1-J



• Molecule 4: Histone H2B type 1-J

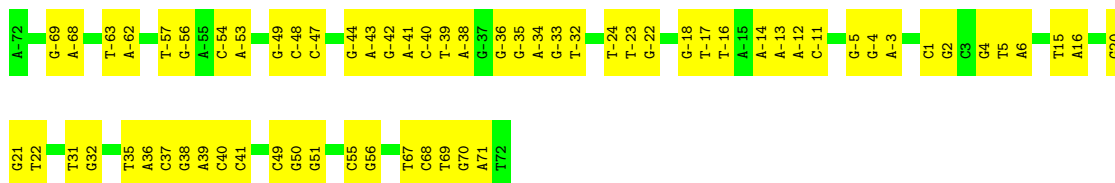


• Molecule 5: DNA (145-MER)

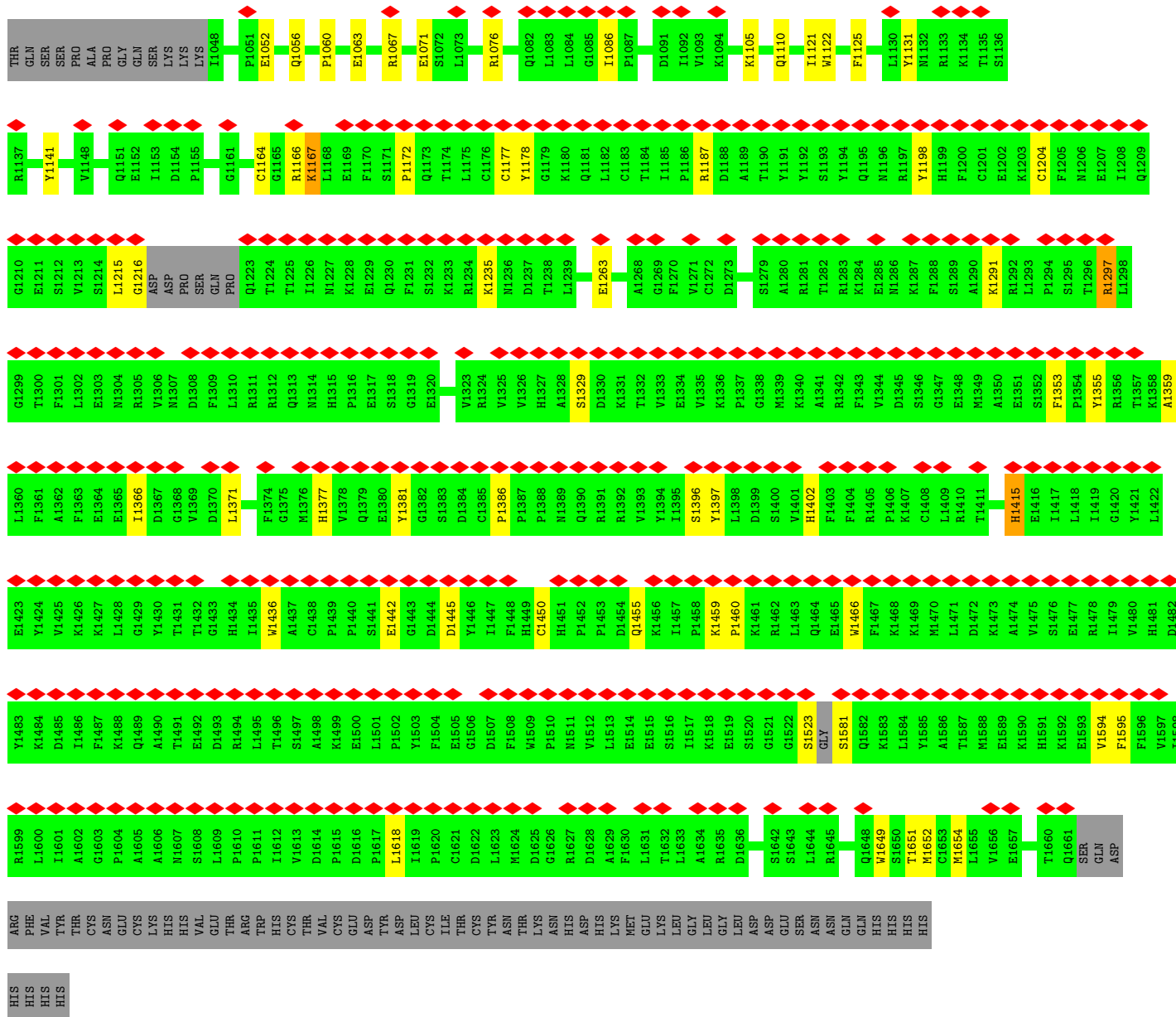
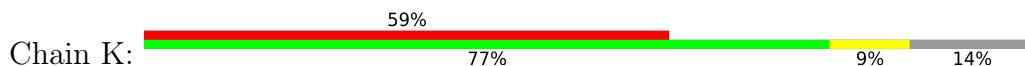


• Molecule 6: DNA (145-MER)





• Molecule 7: Histone acetyltransferase p300



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	25884	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	56	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2300	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.066	Depositor
Minimum map value	-0.035	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0105	Depositor
Map size (\AA)	256.80002, 256.80002, 256.80002	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.07, 1.07, 1.07	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	0.30	0/813	0.42	0/1090
1	E	0.30	0/811	0.41	0/1086
2	B	0.31	0/669	0.44	0/894
2	F	0.33	0/626	0.47	0/837
3	C	0.27	0/806	0.41	0/1089
3	G	0.30	0/820	0.46	0/1107
4	D	0.31	0/757	0.45	0/1015
4	H	0.29	0/736	0.39	0/990
5	I	0.54	0/3308	0.92	0/5099
6	J	0.54	0/3354	0.90	0/5180
7	K	0.50	0/4633	0.69	0/6263
All	All	0.46	0/17333	0.73	0/24650

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	801	207	839	30	0
1	E	800	208	836	37	0
2	B	662	179	709	23	0
2	F	619	166	659	38	0
3	C	796	200	848	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	810	205	866	8	0
4	D	746	191	771	16	0
4	H	725	181	745	14	0
5	I	2952	294	1629	50	0
6	J	2987	305	1630	54	0
7	K	4512	0	4421	42	0
All	All	16410	2136	13953	254	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 8.

All (254) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:119:ILE:HD13	2:F:43:VAL:HG11	1.32	1.10
1:A:118:THR:HG22	2:B:45:ARG:HB3	1.48	0.95
1:A:107:THR:HG23	1:A:123:ASP:HB3	1.47	0.95
1:A:106:ASP:OD2	1:A:131:ARG:NE	2.02	0.92
1:E:120:MET:C	2:F:50:ILE:HD11	1.95	0.86
1:E:118:THR:HG22	2:F:45:ARG:HG2	1.57	0.86
1:E:107:THR:HG23	1:E:123:ASP:HB3	1.54	0.86
1:E:120:MET:CA	2:F:50:ILE:HD11	2.12	0.80
1:A:106:ASP:OD2	1:A:131:ARG:NH2	2.14	0.79
1:A:106:ASP:OD2	1:A:131:ARG:CZ	2.33	0.76
1:E:106:ASP:OD1	1:E:131:ARG:NE	2.20	0.75
1:A:107:THR:CG2	1:A:123:ASP:HB3	2.18	0.74
1:E:121:PRO:HG2	2:F:49:LEU:HB2	1.70	0.73
6:J:-24:DT:H2'	6:J:-23:DT:H71	1.73	0.70
7:K:1366:ILE:HD12	7:K:1371:LEU:HD23	1.75	0.69
5:I:-70:DC:H2'	5:I:-69:DA:C8	2.27	0.69
1:E:106:ASP:OD1	1:E:131:ARG:NH2	2.26	0.68
7:K:1297:ARG:HH11	7:K:1297:ARG:CG	2.06	0.68
1:E:107:THR:HG23	1:E:123:ASP:CB	2.24	0.67
7:K:1060:PRO:HA	7:K:1063:GLU:HG2	1.79	0.65
6:J:68:DC:H2'	6:J:69:DT:H71	1.79	0.64
1:A:131:ARG:HA	1:E:131:ARG:HA	1.80	0.64
4:D:119:THR:HG22	7:K:1076:ARG:NH1	2.12	0.63
2:F:47:SER:OG	2:F:48:GLY:N	2.31	0.63
1:A:130:ILE:O	1:E:131:ARG:HG2	1.99	0.63
3:G:54:VAL:HG23	4:H:110:ALA:HB1	1.79	0.63
1:A:119:ILE:CD1	2:B:43:VAL:HG11	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:63:GLU:O	2:B:67:ARG:NH1	2.32	0.62
3:G:55:LEU:HD11	4:H:70:PHE:HB2	1.80	0.62
5:I:16:DA:H1'	5:I:17:DA:N7	2.14	0.61
5:I:35:DC:H2''	5:I:36:DC:C5	2.35	0.61
6:J:21:DG:H2'	6:J:22:DT:H71	1.82	0.61
7:K:1172:PRO:HG3	7:K:1187:ARG:HD3	1.81	0.61
1:E:118:THR:HA	2:F:45:ARG:O	2.01	0.61
1:E:120:MET:C	2:F:50:ILE:CD1	2.68	0.60
4:H:36:SER:OG	4:H:63:ASN:ND2	2.35	0.60
4:D:119:THR:CG2	7:K:1076:ARG:NH1	2.65	0.60
2:F:92:ARG:NH2	4:H:100:LEU:O	2.35	0.59
1:A:61:LEU:O	2:B:36:ARG:NH2	2.34	0.59
1:A:108:ASN:HB2	2:B:43:VAL:HG22	1.84	0.59
4:D:115:THR:OG1	7:K:1067:ARG:HG2	2.02	0.59
2:B:51:TYR:O	2:B:55:ARG:NH1	2.36	0.59
2:B:92:ARG:NH2	4:D:100:LEU:O	2.36	0.59
1:E:108:ASN:HB2	2:F:43:VAL:HG12	1.85	0.59
7:K:1172:PRO:CG	7:K:1187:ARG:HD3	2.32	0.58
5:I:26:DG:H2''	5:I:27:DG:C8	2.39	0.58
1:A:119:ILE:HD13	2:B:43:VAL:HG21	1.83	0.58
7:K:1105:LYS:HG2	7:K:1110:GLN:OE1	2.04	0.58
1:E:118:THR:HG22	2:F:45:ARG:CG	2.32	0.57
1:A:119:ILE:HD13	2:B:43:VAL:HG11	1.86	0.57
4:D:86:ARG:NH1	4:D:93:GLU:OE2	2.22	0.57
1:A:107:THR:HG23	1:A:123:ASP:CB	2.29	0.57
1:A:118:THR:CG2	2:B:45:ARG:HE	2.18	0.57
5:I:55:DT:H2''	5:I:56:DC:C5	2.40	0.56
1:E:119:ILE:HD13	2:F:43:VAL:CG1	2.21	0.56
2:B:68:ASP:OD2	2:B:92:ARG:NH1	2.39	0.55
7:K:1235:LYS:HD2	7:K:1235:LYS:N	2.21	0.55
1:A:68:GLN:HE21	1:A:72:ARG:HH21	1.55	0.55
1:E:106:ASP:CG	1:E:131:ARG:HH21	2.10	0.55
1:E:121:PRO:CG	2:F:49:LEU:HB2	2.36	0.55
2:F:68:ASP:OD2	2:F:92:ARG:NH1	2.40	0.55
2:F:78:ARG:NH1	2:F:82:THR:OG1	2.40	0.55
6:J:-69:DG:H2''	6:J:-68:DA:C8	2.41	0.55
6:J:-17:DT:H4'	6:J:-16:DT:OP1	2.07	0.54
4:D:31:ARG:NH1	6:J:51:DG:OP1	2.41	0.54
5:I:-57:DC:H2''	5:I:-56:DC:C5	2.42	0.54
1:A:83:ARG:HB2	2:B:80:THR:HG22	1.90	0.54
3:C:90:ASP:OD1	3:C:91:GLU:N	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD12	2:B:37:LEU:HD23	1.90	0.54
6:J:49:DC:H2''	6:J:50:DG:C8	2.43	0.54
1:E:106:ASP:OD1	1:E:131:ARG:CZ	2.57	0.53
5:I:46:DA:H2''	5:I:47:DG:N7	2.24	0.52
5:I:49:DC:H2''	5:I:50:DA:C8	2.44	0.52
6:J:-57:DT:H2''	6:J:-56:DG:C8	2.44	0.52
2:B:64:ASN:O	2:B:93:GLN:NE2	2.42	0.52
4:D:112:SER:HA	7:K:1067:ARG:HD2	1.91	0.52
1:A:118:THR:HA	2:B:45:ARG:O	2.09	0.52
3:C:55:LEU:HD11	4:D:70:PHE:HB2	1.92	0.52
5:I:-27:DC:H2''	5:I:-26:DT:H71	1.91	0.52
6:J:67:DT:H2''	6:J:68:DC:C6	2.45	0.52
6:J:-63:DT:H2''	6:J:-62:DA:H8	1.75	0.51
1:E:119:ILE:HG21	2:F:43:VAL:CG1	2.40	0.51
6:J:-34:DA:H2''	6:J:-33:DG:C8	2.45	0.51
7:K:1198:TYR:CD2	7:K:1215:LEU:HD13	2.44	0.51
7:K:1359:ALA:HB2	7:K:1377:HIS:CD2	2.45	0.51
1:A:103:LEU:CD1	1:A:128:ARG:HG3	2.39	0.51
5:I:68:DT:H2''	5:I:69:DC:C5	2.46	0.51
1:A:103:LEU:HD13	1:A:128:ARG:HG3	1.93	0.51
7:K:1402:HIS:HD2	7:K:1455:GLN:HB2	1.75	0.51
1:E:61:LEU:HD12	2:F:37:LEU:HD23	1.93	0.51
1:E:121:PRO:N	2:F:50:ILE:HD13	2.26	0.51
4:H:37:TYR:H	4:H:63:ASN:ND2	2.09	0.51
6:J:-4:DG:H2''	6:J:-3:DA:C8	2.46	0.50
5:I:-26:DT:H2''	5:I:-25:DA:C8	2.46	0.50
6:J:-5:DG:H2''	6:J:-4:DG:C8	2.47	0.50
6:J:-18:DG:H2''	6:J:-17:DT:C5	2.47	0.50
1:E:120:MET:HA	2:F:50:ILE:HD11	1.91	0.49
5:I:-54:DA:H2''	5:I:-53:DG:C8	2.47	0.49
7:K:1297:ARG:HH11	7:K:1297:ARG:HG3	1.75	0.49
1:E:61:LEU:O	2:F:36:ARG:NH2	2.39	0.49
5:I:-14:DA:H2''	5:I:-13:DA:H8	1.77	0.49
6:J:-33:DG:H2'	6:J:-32:DT:H71	1.94	0.49
1:A:71:VAL:HG13	2:B:66:ILE:HD11	1.95	0.49
4:D:119:THR:HG22	7:K:1076:ARG:HH11	1.74	0.49
2:B:98:TYR:CE2	3:G:100:VAL:HG21	2.48	0.49
3:G:29:ARG:NH1	4:H:36:SER:O	2.46	0.49
5:I:-6:DT:H2''	5:I:-5:DA:N7	2.27	0.49
1:E:51:ILE:HD11	2:F:43:VAL:O	2.13	0.48
5:I:4:DC:H2''	5:I:5:DC:C5	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:K:1052:GLU:O	7:K:1056:GLN:HG2	2.13	0.48
2:F:38:ALA:HB1	2:F:43:VAL:HG21	1.96	0.48
6:J:39:DA:H2''	6:J:40:DC:H5'	1.95	0.48
5:I:7:DC:H5''	5:I:7:DC:H6	1.78	0.48
6:J:-54:DC:H2''	6:J:-53:DA:C8	2.49	0.48
1:A:83:ARG:O	2:B:80:THR:HA	2.14	0.47
1:E:119:ILE:HG21	2:F:43:VAL:HG12	1.96	0.47
1:E:121:PRO:N	2:F:50:ILE:CD1	2.76	0.47
5:I:-47:DT:H2''	5:I:-46:DC:C5	2.49	0.47
6:J:-44:DG:H2''	6:J:-43:DA:H8	1.79	0.47
3:G:92:GLU:OE2	4:H:105:GLU:HB2	2.13	0.47
6:J:21:DG:H2''	6:J:22:DT:O5'	2.14	0.47
5:I:-14:DA:H2''	5:I:-13:DA:C8	2.49	0.47
6:J:-23:DT:C2	6:J:-22:DG:C8	3.03	0.47
6:J:-17:DT:H2''	6:J:-16:DT:O5'	2.14	0.47
5:I:-36:DT:H2''	5:I:-35:DA:N7	2.29	0.47
5:I:64:DT:H2''	5:I:65:DA:C8	2.50	0.47
7:K:1297:ARG:HH11	7:K:1297:ARG:HG2	1.80	0.47
6:J:39:DA:C4	6:J:40:DC:C5	3.03	0.47
2:B:35:ARG:NH2	2:B:51:TYR:OH	2.47	0.47
6:J:-43:DA:H2''	6:J:-42:DG:H8	1.80	0.47
7:K:1071:GLU:HA	7:K:1141:TYR:CD2	2.49	0.47
7:K:1355:TYR:HB3	7:K:1381:TYR:CE2	2.50	0.47
5:I:-68:DG:H2''	5:I:-67:DA:C8	2.50	0.46
5:I:42:DC:H2''	5:I:43:DT:O5'	2.14	0.46
7:K:1353:PHE:CD2	7:K:1386:PRO:HG2	2.50	0.46
4:H:111:VAL:O	4:H:115:THR:HG23	2.15	0.46
5:I:-8:DC:H2''	5:I:-7:DG:C8	2.51	0.46
6:J:31:DT:H2''	6:J:32:DG:C8	2.50	0.46
4:D:38:SER:HA	4:D:59:MET:HE1	1.96	0.46
2:F:75:HIS:HB2	4:H:96:THR:HG21	1.98	0.46
6:J:35:DT:H2''	6:J:36:DA:C8	2.51	0.46
6:J:-34:DA:H2''	6:J:-33:DG:H8	1.81	0.46
4:D:119:THR:CG2	7:K:1076:ARG:HH12	2.27	0.46
7:K:1366:ILE:CD1	7:K:1371:LEU:HD23	2.44	0.46
1:E:83:ARG:O	2:F:80:THR:HA	2.16	0.45
3:G:65:LEU:HB2	3:G:86:ALA:HB1	1.98	0.45
7:K:1436:TRP:CD1	7:K:1595:PHE:CE1	3.05	0.45
4:H:106:LEU:HD23	4:H:106:LEU:HA	1.86	0.45
5:I:8:DC:H2''	5:I:9:DG:C8	2.51	0.45
6:J:20:DG:H2''	6:J:21:DG:H8	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:35:ARG:NH1	5:I:8:DC:OP2	2.37	0.45
1:E:100:LEU:HD23	1:E:100:LEU:HA	1.78	0.45
5:I:53:DT:H2''	5:I:54:DG:C8	2.51	0.45
7:K:1415:HIS:HE1	7:K:1466:TRP:CE2	2.34	0.45
5:I:-48:DC:H2''	5:I:-47:DT:C5	2.52	0.45
5:I:19:DC:H2''	5:I:20:DG:H8	1.82	0.45
1:E:68:GLN:HE21	1:E:72:ARG:HH21	1.65	0.45
5:I:17:DA:H1'	5:I:18:DC:H5'	1.99	0.45
7:K:1086:ILE:HG23	7:K:1131:TYR:OH	2.17	0.45
6:J:-48:DC:H2''	6:J:-47:DC:C6	2.52	0.45
1:E:125:GLN:NE2	2:F:53:GLU:OE1	2.50	0.45
6:J:-49:DG:C8	6:J:-49:DG:H5'	2.51	0.45
7:K:1649:TRP:CZ3	7:K:1652:MET:CE	3.00	0.45
6:J:-36:DG:H2''	6:J:-35:DG:C8	2.52	0.44
6:J:-63:DT:H2''	6:J:-62:DA:C8	2.53	0.44
5:I:-38:DC:H2''	5:I:-37:DG:C8	2.53	0.44
1:E:121:PRO:HB3	2:F:53:GLU:HB2	2.00	0.44
5:I:-66:DA:C8	5:I:-66:DA:H5'	2.52	0.44
5:I:16:DA:H1'	5:I:17:DA:C8	2.52	0.44
1:A:72:ARG:HH22	5:I:-23:DC:P	2.41	0.44
6:J:1:DC:H2''	6:J:2:DG:C8	2.53	0.44
7:K:1402:HIS:CD2	7:K:1455:GLN:HB2	2.53	0.44
5:I:-5:DA:N6	6:J:4:DG:O6	2.50	0.44
5:I:-2:DC:H2''	5:I:-1:DG:C8	2.52	0.44
5:I:67:DA:H5'	5:I:67:DA:C8	2.52	0.44
6:J:49:DC:H2''	6:J:50:DG:H8	1.81	0.44
6:J:55:DC:H2''	6:J:56:DG:C8	2.52	0.44
1:E:107:THR:CG2	1:E:123:ASP:HB3	2.37	0.44
6:J:-14:DA:H2''	6:J:-13:DA:H8	1.82	0.43
6:J:40:DC:C2	6:J:41:DC:C5	3.06	0.43
6:J:70:DG:H2''	6:J:71:DA:C8	2.52	0.43
6:J:69:DT:H2''	6:J:70:DG:H8	1.83	0.43
1:A:118:THR:HG22	2:B:45:ARG:CB	2.34	0.43
2:F:37:LEU:HD23	2:F:37:LEU:HA	1.90	0.43
5:I:60:DT:H2''	5:I:61:DA:H8	1.84	0.43
6:J:-39:DT:H2''	6:J:-38:DA:C8	2.53	0.43
7:K:1121:ILE:HG22	7:K:1125:PHE:CE2	2.53	0.43
1:A:62:ILE:O	1:A:93:GLN:NE2	2.51	0.43
1:A:113:HIS:CG	1:E:126:LEU:HD22	2.54	0.43
4:D:46:LYS:HA	4:D:46:LYS:HD3	1.79	0.43
1:E:125:GLN:NE2	2:F:53:GLU:CD	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:-58:DG:H2''	5:I:-57:DC:C5	2.54	0.42
7:K:1122:TRP:CD2	7:K:1167:LYS:HD2	2.54	0.42
4:H:38:SER:HA	4:H:59:MET:HE1	2.01	0.42
5:I:4:DC:H2''	5:I:5:DC:C6	2.54	0.42
7:K:1177:CYS:SG	7:K:1204:CYS:SG	3.16	0.42
6:J:-41:DA:H1'	6:J:-40:DC:H5'	2.02	0.42
6:J:55:DC:H2''	6:J:56:DG:N7	2.34	0.42
1:A:100:LEU:HA	1:A:100:LEU:HD23	1.76	0.42
3:C:54:VAL:HG21	4:D:98:VAL:HG21	2.00	0.42
3:C:100:VAL:HG21	2:F:98:TYR:CE2	2.54	0.42
4:D:107:ALA:O	4:D:111:VAL:HG23	2.19	0.42
2:F:62:LEU:HD23	2:F:62:LEU:HA	1.77	0.42
4:H:108:LYS:HB2	4:H:108:LYS:HE3	1.62	0.42
6:J:-12:DA:C4	6:J:-11:DC:C5	3.07	0.42
6:J:37:DC:H2''	6:J:38:DG:C8	2.54	0.42
1:E:71:VAL:HG13	2:F:66:ILE:HD11	2.00	0.42
6:J:38:DG:H2''	6:J:39:DA:C8	2.54	0.42
7:K:1523:SER:OG	7:K:1581:SER:N	2.52	0.42
2:F:46:ILE:HG22	2:F:47:SER:O	2.20	0.42
5:I:22:DC:H2''	5:I:23:DA:H8	1.84	0.42
7:K:1445:ASP:OD2	7:K:1450:CYS:N	2.44	0.42
7:K:1459:LYS:HB3	7:K:1460:PRO:CD	2.50	0.42
3:C:23:LEU:HD23	3:C:23:LEU:HA	1.86	0.41
3:G:42:ARG:O	4:H:88:THR:HA	2.20	0.41
5:I:-6:DT:H2''	5:I:-5:DA:C8	2.55	0.41
6:J:15:DT:H2''	6:J:16:DA:C8	2.55	0.41
5:I:5:DC:H2''	5:I:6:DC:C5	2.56	0.41
6:J:-14:DA:H2''	6:J:-13:DA:C8	2.55	0.41
7:K:1442:GLU:H	7:K:1442:GLU:CD	2.24	0.41
7:K:1651:THR:HA	7:K:1654:MET:HE3	2.03	0.41
2:B:37:LEU:HD23	2:B:37:LEU:HA	1.89	0.41
2:B:38:ALA:HB3	2:B:46:ILE:HD11	2.02	0.41
2:B:62:LEU:HD23	2:B:62:LEU:HA	1.79	0.41
5:I:61:DA:H2'	5:I:62:DT:H71	2.03	0.41
7:K:1122:TRP:CG	7:K:1167:LYS:HD2	2.55	0.41
5:I:10:DC:H2''	5:I:11:DG:C8	2.56	0.41
5:I:45:DC:H2''	5:I:46:DA:C8	2.55	0.41
7:K:1436:TRP:CD1	7:K:1594:VAL:HG12	2.56	0.41
3:C:63:LEU:HD13	4:D:45:LEU:HB2	2.02	0.41
6:J:-35:DG:H2''	6:J:-34:DA:H8	1.86	0.41
6:J:68:DC:H2'	6:J:69:DT:C6	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:J:69:DT:H2''	6:J:70:DG:C8	2.56	0.41
5:I:7:DC:H5''	5:I:7:DC:C6	2.56	0.41
7:K:1172:PRO:HG2	7:K:1187:ARG:HD3	2.00	0.41
7:K:1329:SER:HA	7:K:1618:LEU:HD23	2.03	0.41
1:A:68:GLN:NE2	1:A:72:ARG:HH21	2.19	0.40
3:C:62:ILE:HD11	3:C:93:LEU:HD22	2.03	0.40
4:D:100:LEU:HD23	4:D:100:LEU:HA	1.87	0.40
2:F:73:THR:OG1	2:F:85:ASP:OD2	2.31	0.40
1:A:103:LEU:HD21	1:A:124:ILE:HG23	2.04	0.40
5:I:1:DT:H2''	5:I:2:DG:C8	2.55	0.40
7:K:1178:TYR:OH	7:K:1216:GLY:O	2.39	0.40
3:C:33:LEU:HD23	3:C:33:LEU:HA	1.88	0.40
2:F:77:LYS:HE3	4:H:92:ARG:NH1	2.36	0.40
3:G:34:LEU:HA	3:G:34:LEU:HD23	1.88	0.40
5:I:-25:DA:H2''	5:I:-24:DG:H8	1.85	0.40
6:J:-36:DG:H2''	6:J:-35:DG:N7	2.36	0.40
6:J:5:DT:H2''	6:J:6:DA:N7	2.36	0.40
1:E:66:PRO:HD3	5:I:17:DA:H5''	2.04	0.40
5:I:-66:DA:H5'	5:I:-66:DA:H8	1.85	0.40
6:J:-23:DT:C2	6:J:-22:DG:N7	2.89	0.40
2:F:30:THR:HG21	6:J:-13:DA:H5''	2.03	0.40
7:K:1164:CYS:SG	7:K:1166:ARG:HG3	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	95/135 (70%)	94 (99%)	1 (1%)	0	100	100
1	E	95/135 (70%)	93 (98%)	2 (2%)	0	100	100
2	B	81/106 (76%)	78 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	76/106 (72%)	73 (96%)	3 (4%)	0	100	100
3	C	101/133 (76%)	101 (100%)	0	0	100	100
3	G	103/133 (77%)	103 (100%)	0	0	100	100
4	D	93/129 (72%)	92 (99%)	1 (1%)	0	100	100
4	H	91/129 (70%)	90 (99%)	1 (1%)	0	100	100
7	K	545/638 (85%)	530 (97%)	15 (3%)	0	100	100
All	All	1280/1644 (78%)	1254 (98%)	26 (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	85/110 (77%)	85 (100%)	0	100	100
1	E	84/110 (76%)	84 (100%)	0	100	100
2	B	68/81 (84%)	68 (100%)	0	100	100
2	F	63/81 (78%)	63 (100%)	0	100	100
3	C	82/102 (80%)	82 (100%)	0	100	100
3	G	83/102 (81%)	83 (100%)	0	100	100
4	D	81/107 (76%)	81 (100%)	0	100	100
4	H	79/107 (74%)	79 (100%)	0	100	100
7	K	501/583 (86%)	494 (99%)	7 (1%)	67	80
All	All	1126/1383 (81%)	1119 (99%)	7 (1%)	86	91

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
7	K	1167	LYS
7	K	1263	GLU

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Mol	Chain	Res	Type
7	K	1291	LYS
7	K	1297	ARG
7	K	1396	SER
7	K	1397	TYR
7	K	1415	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
2	B	25	ASN
2	B	75	HIS
4	D	84	ASN
1	E	39	HIS
1	E	68	GLN
3	G	24	GLN
3	G	73	ASN
4	H	63	ASN
7	K	1256	GLN
7	K	1402	HIS
7	K	1415	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

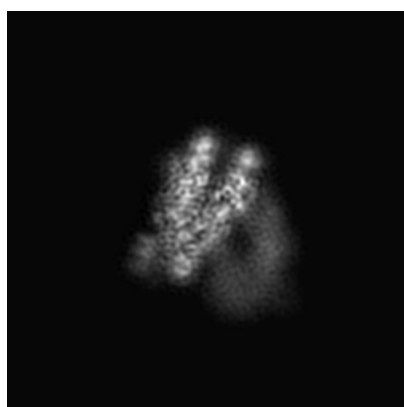
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32373. These allow visual inspection of the internal detail of the map and identification of artifacts.

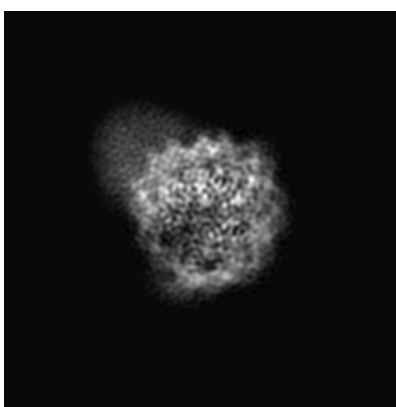
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

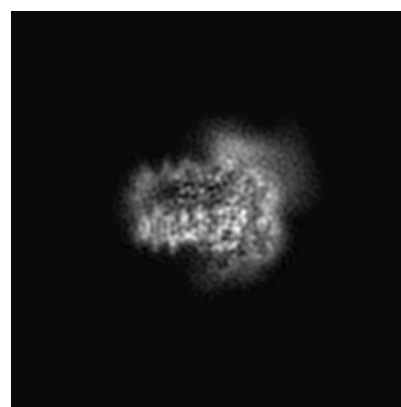
6.1.1 Primary map



X



Y

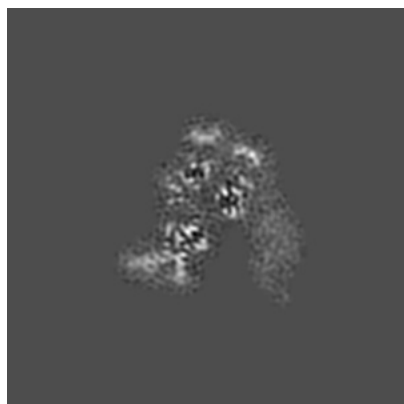


Z

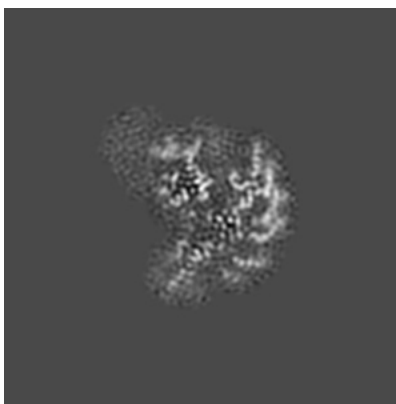
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

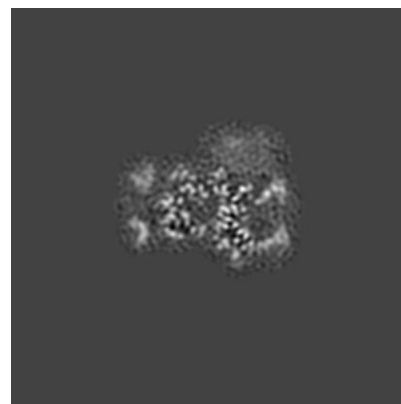
6.2.1 Primary map



X Index: 120



Y Index: 120

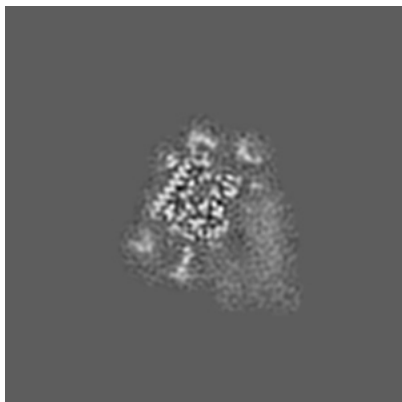


Z Index: 120

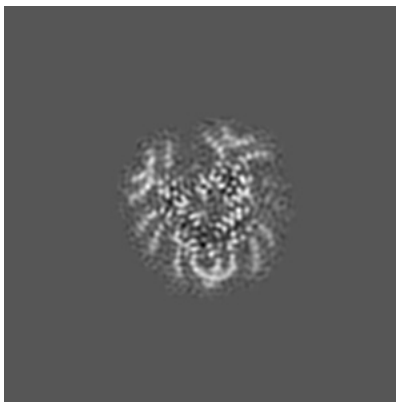
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

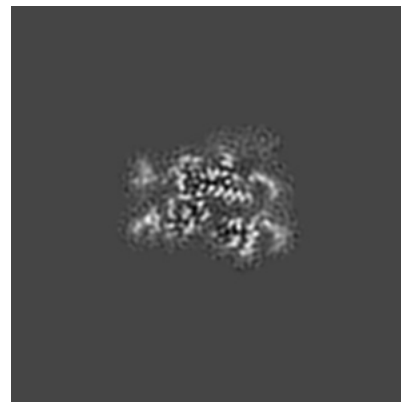
6.3.1 Primary map



X Index: 134



Y Index: 109



Z Index: 129

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

6.4 Orthogonal surface views [i](#)

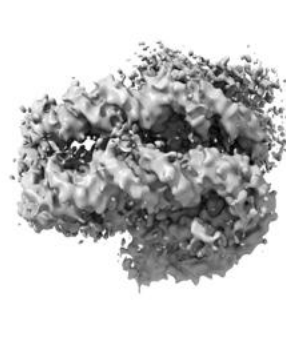
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0105. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

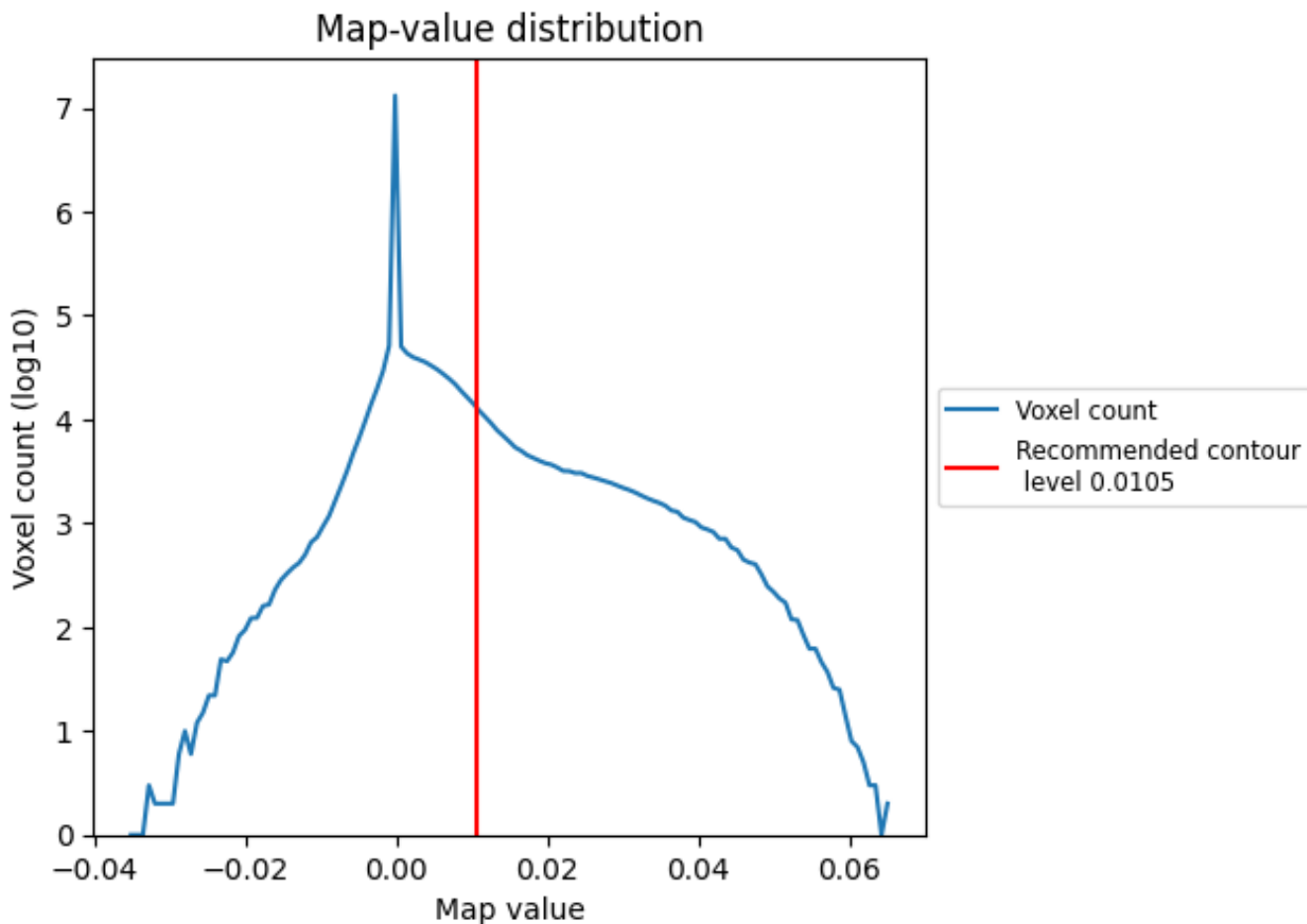
6.5 Mask visualisation

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

7 Map analysis [i](#)

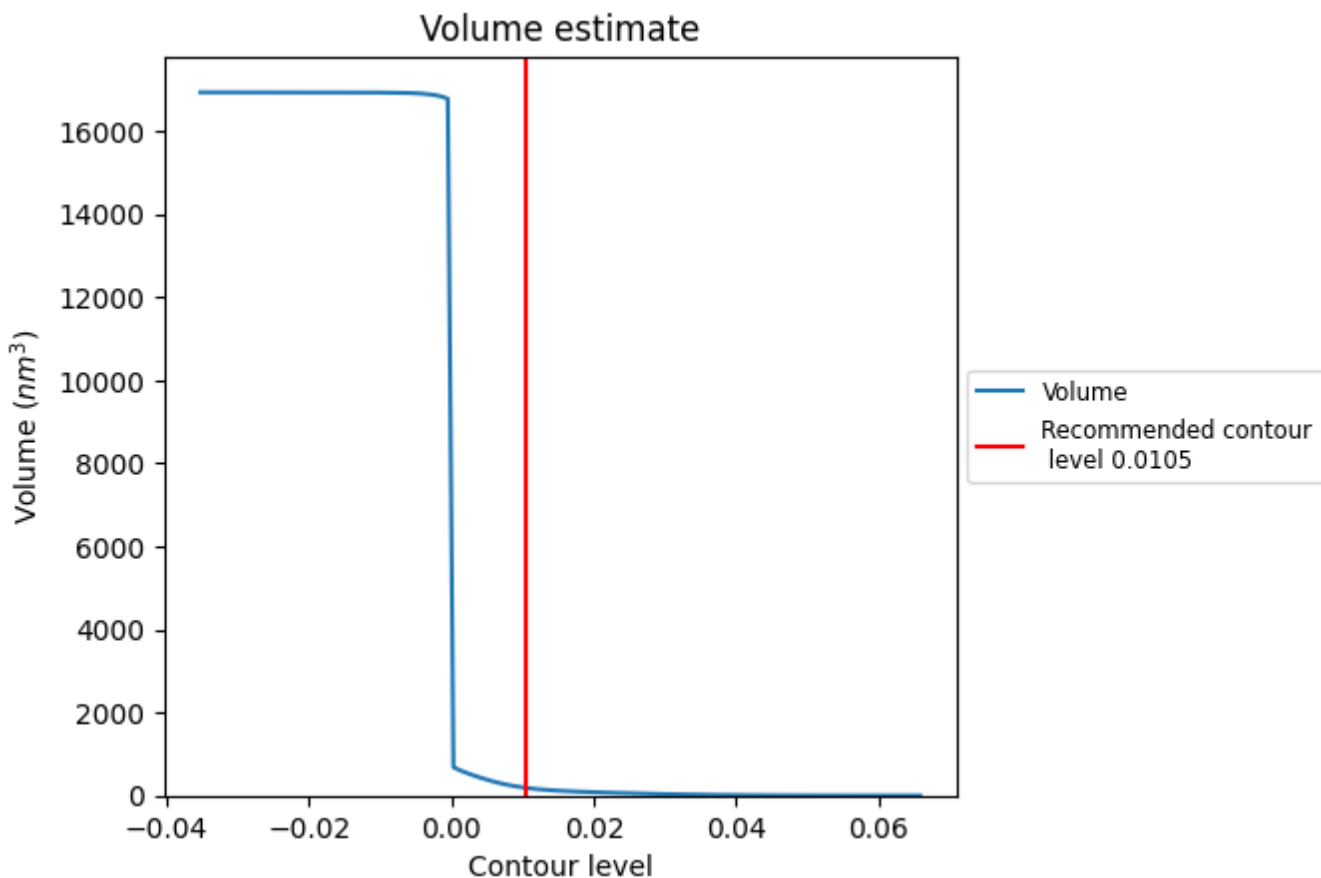
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

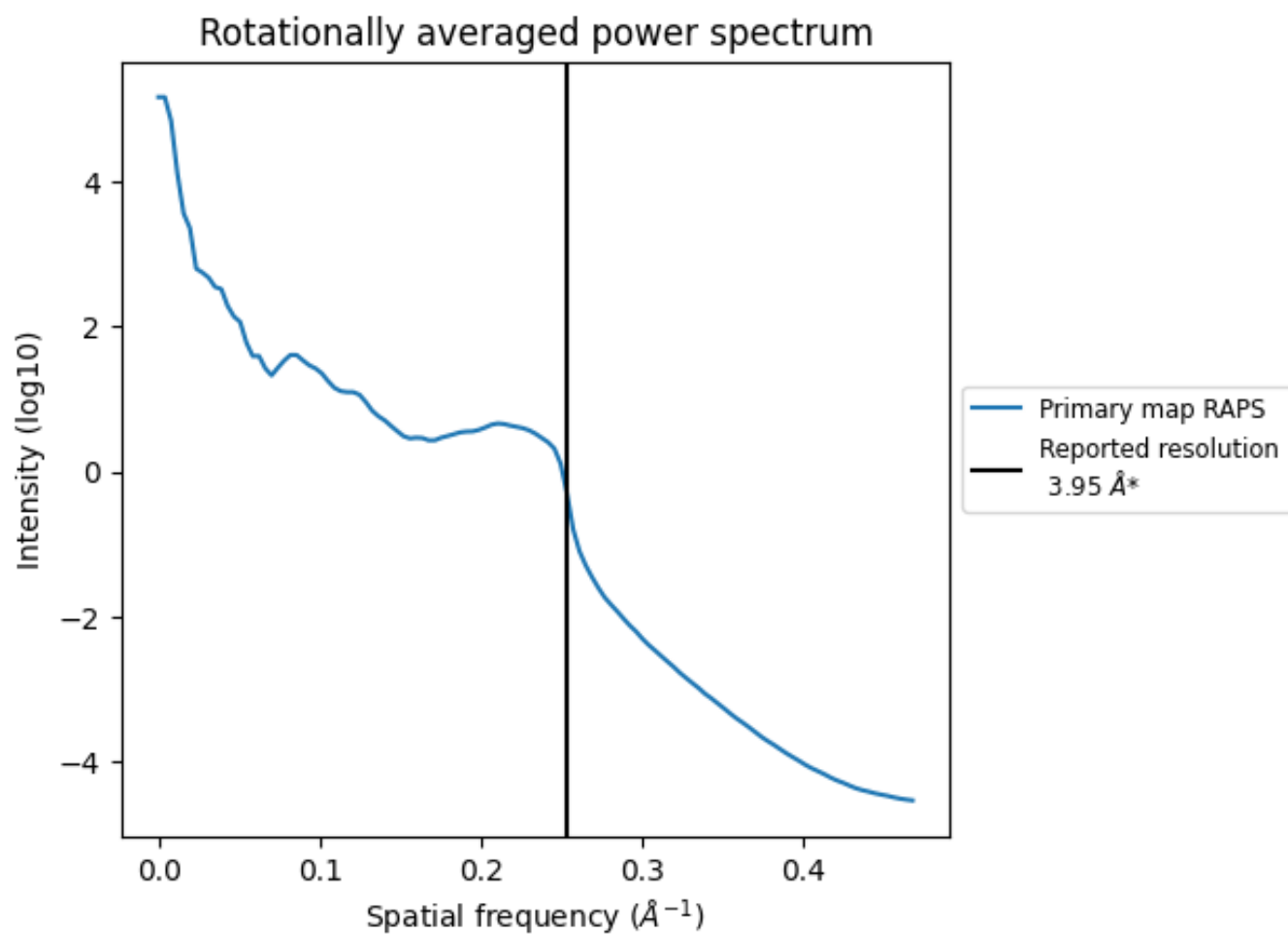
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 186 nm³; this corresponds to an approximate mass of 168 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

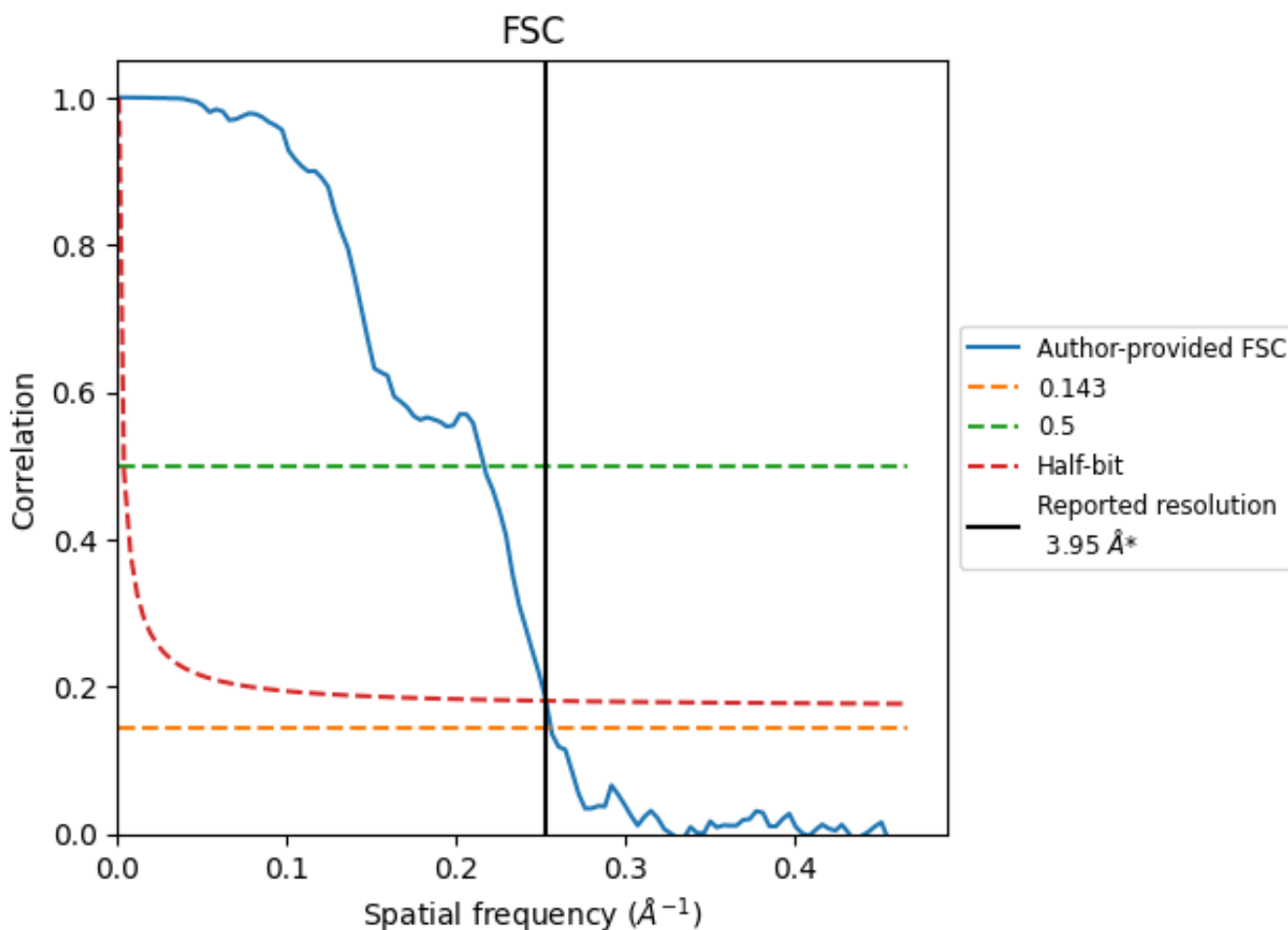


*Reported resolution corresponds to spatial frequency of 0.253 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.253 Å⁻¹

8.2 Resolution estimates [i](#)

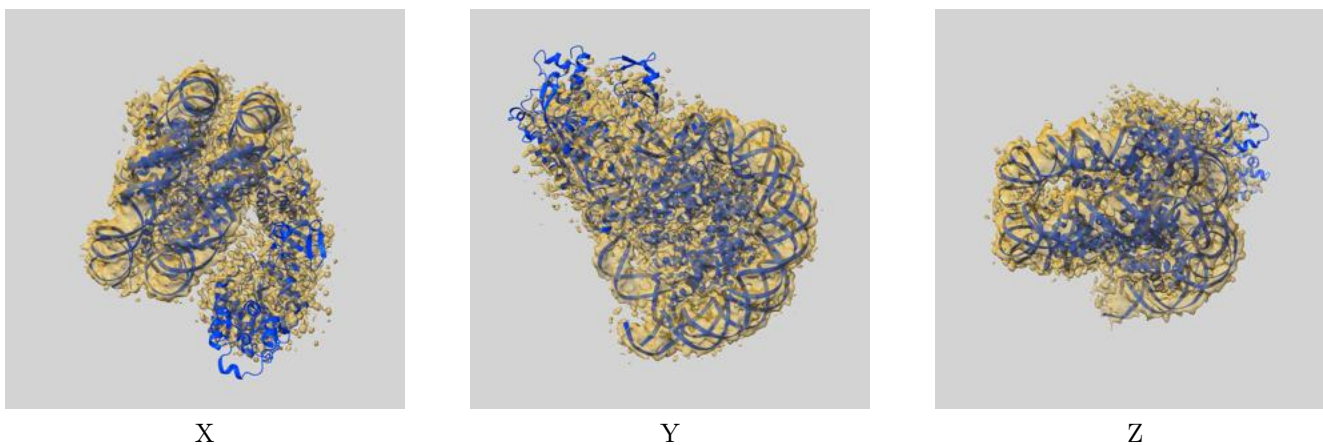
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.95	-	-
Author-provided FSC curve	3.90	4.61	3.95
Unmasked-calculated*	-	-	-

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

9 Map-model fit [i](#)

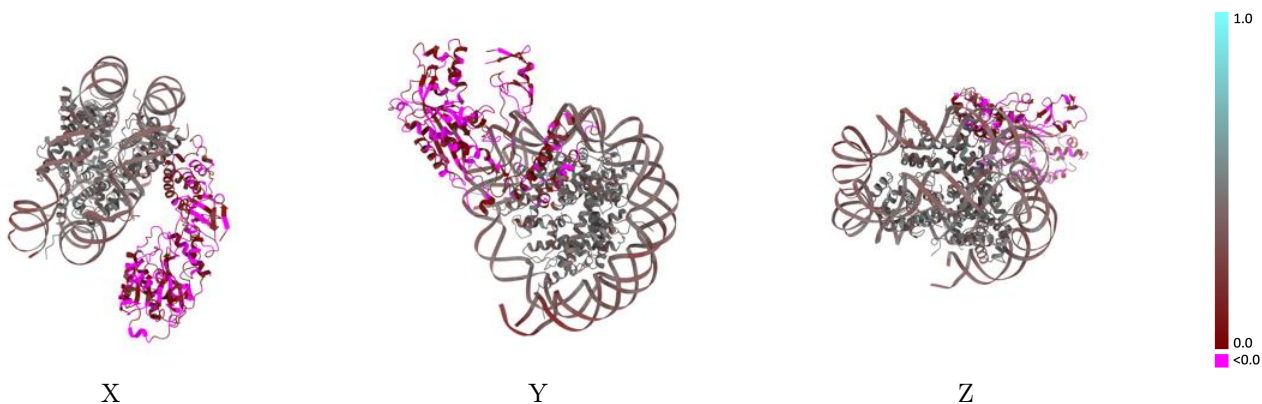
This section contains information regarding the fit between EMDB map EMD-32373 and PDB model 7W9V. Per-residue inclusion information can be found in section [3](#) on page [7](#).

9.1 Map-model overlay [i](#)



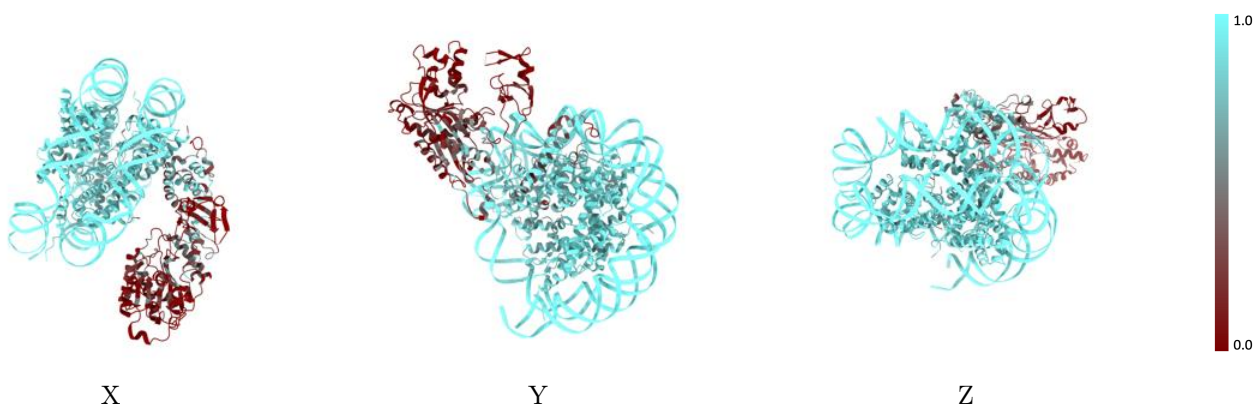
The images above show the 3D surface view of the map at the recommended contour level 0.0105 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



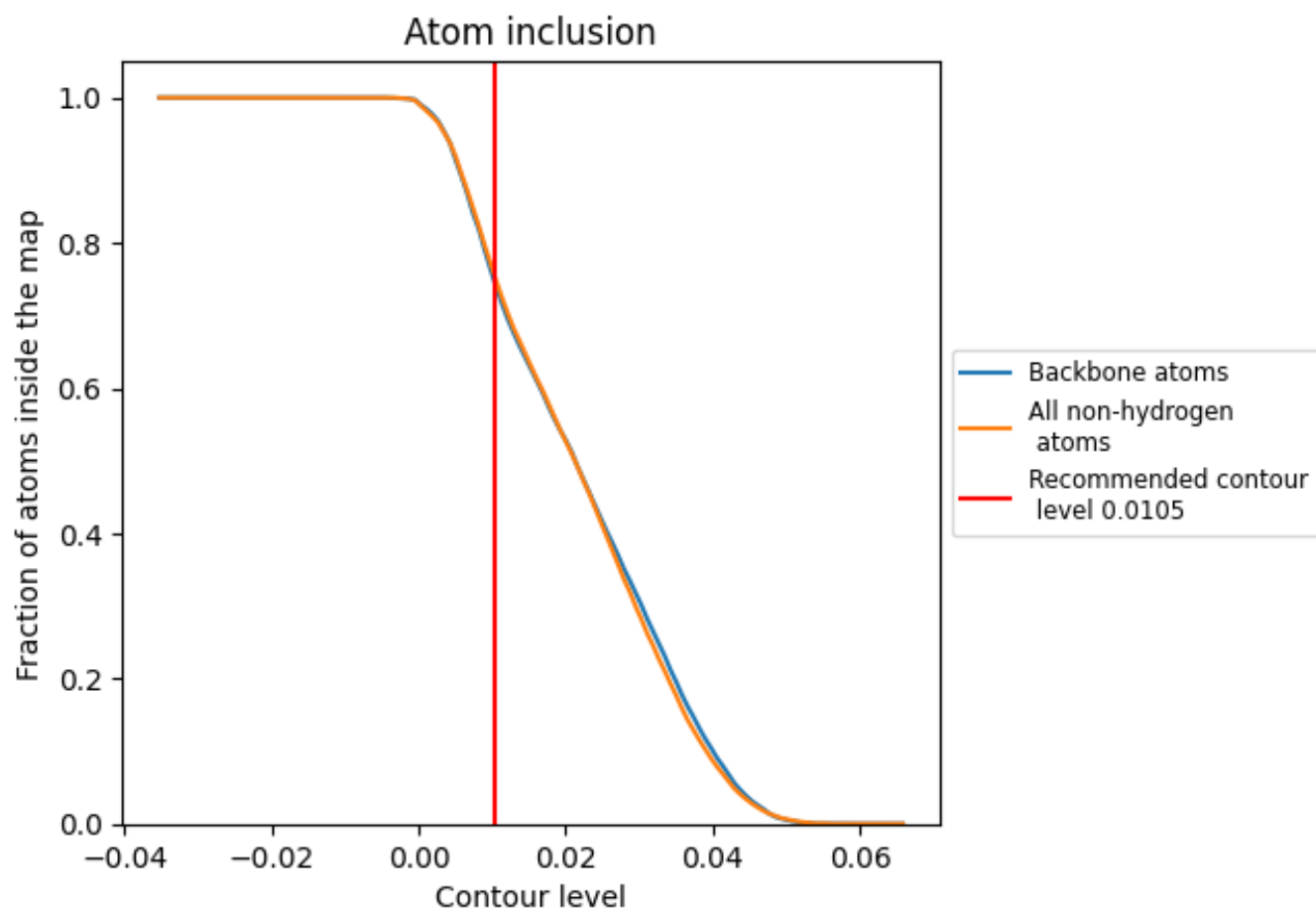
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0105).

























9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 75% 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 (0.0105) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7501	 0.3070
A	 0.9065	 0.4520
B	 0.9057	 0.4560
C	 0.9170	 0.4440
D	 0.9175	 0.4330
E	 0.9103	 0.4400
F	 0.9479	 0.4700
G	 0.9057	 0.4520
H	 0.9463	 0.4490
I	 0.9936	 0.3650
J	 0.9926	 0.3670
K	 0.2959	 0.0400

