



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 03:17 PM UTC

PDB ID : 6KIZ / pdb_00006kiz
EMDB ID : EMD-0695
Title : Cryo-EM structure of human MLL1-NCP complex, binding mode2
Authors : Huang, J.; Xue, H.; Yao, T.
Deposited on : 2019-07-20
Resolution : 4.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

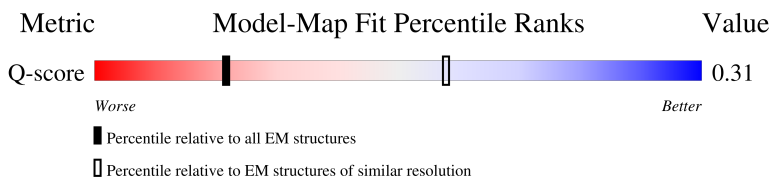
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.50 Å.

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



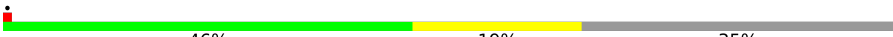
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2937 (4.00 - 5.00)

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	 64% 7% 28%
1	E	135	 67% . 28%
2	B	102	 69% 12% 20%
2	F	102	 75% . 24%

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Mol	Chain	Length	Quality of chain
3	C	129	 79%17%
3	G	129	 80%19%
4	D	122	 73%23%
4	H	122	 72%6%22%
5	I	145	 87%13%
6	J	145	 85%15%
7	N	538	 46%19%35%
8	K	216	 20%64%19%17%
9	R	334	 66%24%10%
10	T	534	 25%8%67%

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 19889 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Histone H3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	97	Total	C	N	O	S	0	0
			802	506	155	138	3		
1	E	97	Total	C	N	O	S	0	0
			802	506	155	138	3		

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	82	Total	C	N	O	S	0	0
			657	416	128	112	1		
2	F	78	Total	C	N	O	S	0	0
			622	393	120	108	1		

- Molecule 3 is a protein called Histone H2A.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	107	Total	C	N	O	0	0
			825	519	163	143		
3	G	105	Total	C	N	O	0	0
			809	510	158	141		

- Molecule 4 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	94	Total	C	N	O	S	0	0
			738	463	134	138	3		
4	H	95	Total	C	N	O	S	0	0
			742	466	133	140	3		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	29	THR	SER	engineered mutation	UNP P02281

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Chain	Residue	Modelled	Actual	Comment	Reference
D	117	CYS	LYS	engineered mutation	UNP P02281
H	29	THR	SER	engineered mutation	UNP P02281
H	117	CYS	LYS	engineered mutation	UNP P02281

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	145	Total	C	N	O	P	0	0
			2955	1403	538	869	145		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	145	Total	C	N	O	P	0	0
			2990	1415	559	871	145		

- Molecule 7 is a protein called Retinoblastoma-binding protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	N	351	Total	C	N	O	S	0	0
			2759	1738	477	529	15		

- Molecule 8 is a protein called Histone-lysine N-methyltransferase 2A.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	K	179	Total	C	N	O	S	0	0
			1436	901	268	252	15		

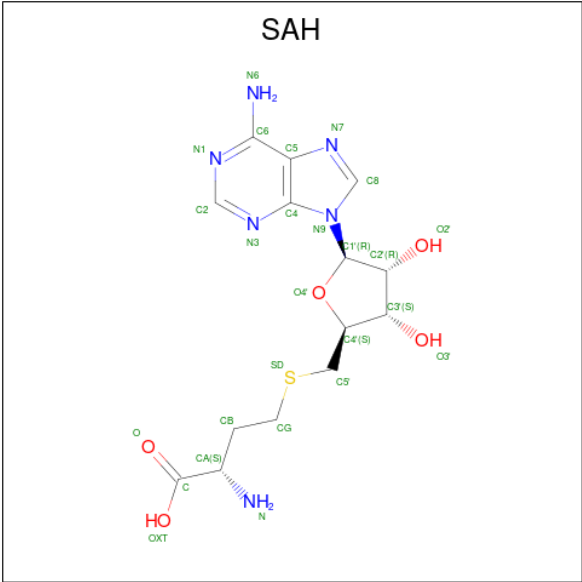
- Molecule 9 is a protein called WD repeat-containing protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	R	300	Total	C	N	O	S	0	0
			2326	1485	388	444	9		

- Molecule 10 is a protein called Set1/Ash2 histone methyltransferase complex subunit ASH2.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	T	176	Total	C	N	O	S	0	0
			1399	906	234	253	6		

- Molecule 11 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: C₁₄H₂₀N₆O₅S).



Mol	Chain	Residues	Atoms					AltConf
11	K	1	Total	C	N	O	S	0
			26	14	6	5	1	

- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn).

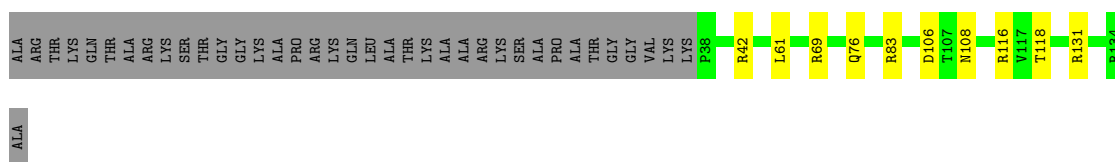
Mol	Chain	Residues	Atoms		AltConf
12	K	1	Total	Zn	0
			1	1	

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.

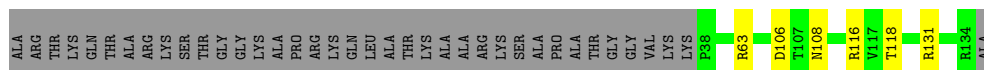
• Molecule 1: Histone H3

Chain A: 



• Molecule 1: Histone H3

Chain E: 



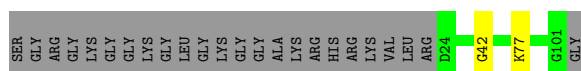
• Molecule 2: Histone H4

Chain B: 



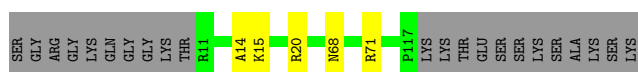
• Molecule 2: Histone H4

Chain F: 



• Molecule 3: Histone H2A

Chain C: 



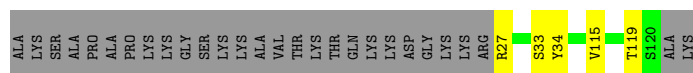
• Molecule 3: Histone H2A

Chain G:  80% 19%



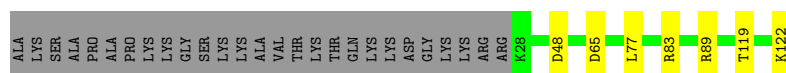
- Molecule 4: Histone H2B 1.1

Chain D:  73% 23%



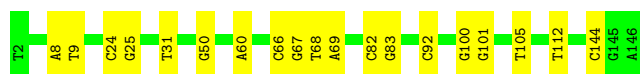
- Molecule 4: Histone H2B 1.1

Chain H:  72% 22%



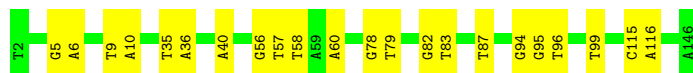
- Molecule 5: DNA (145-MER)

Chain I:  87% 13%



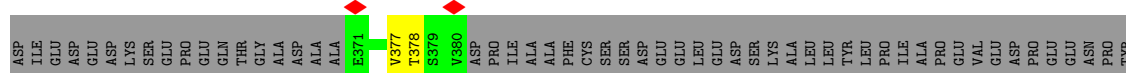
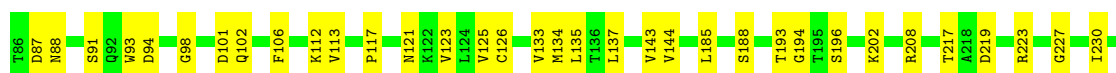
- Molecule 6: DNA (145-MER)

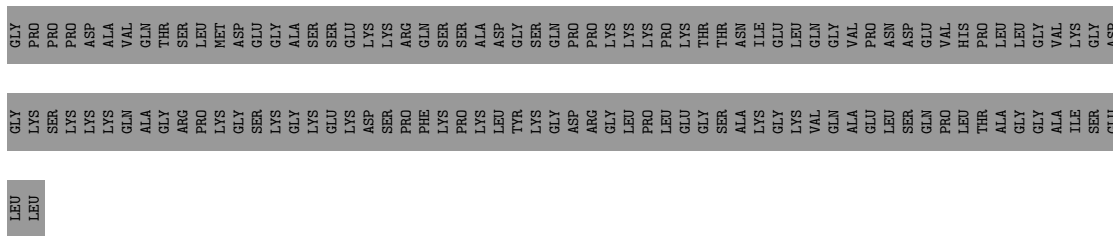
Chain J:  85% 15%



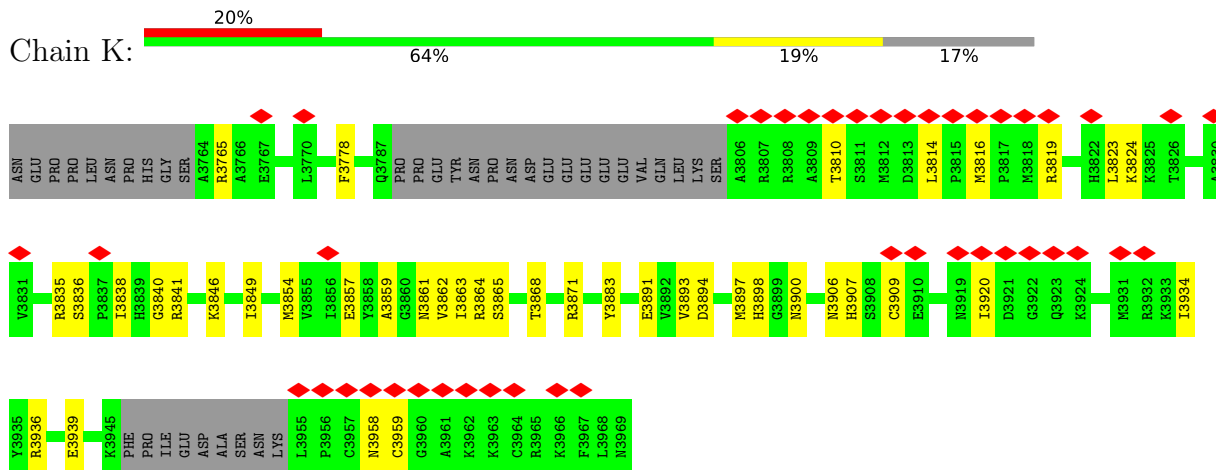
- Molecule 7: Retinoblastoma-binding protein 5

Chain N:  46% 19% 35%

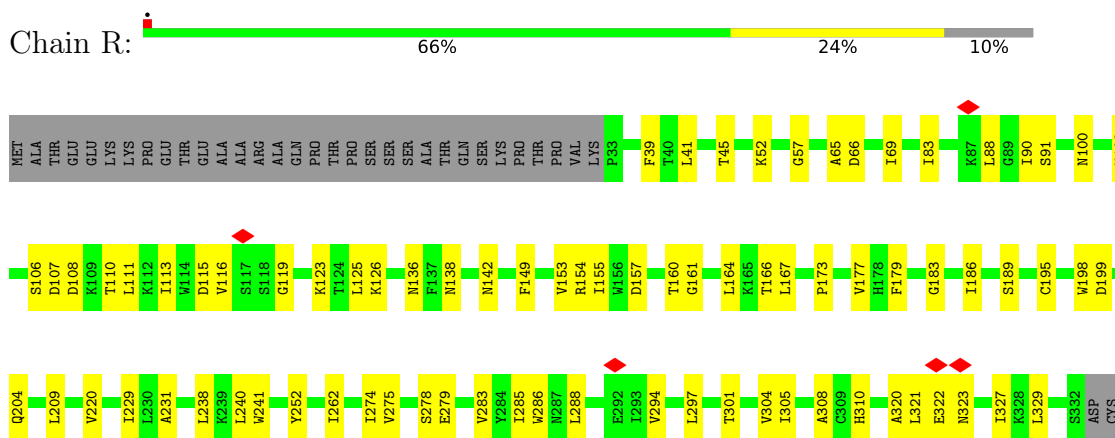




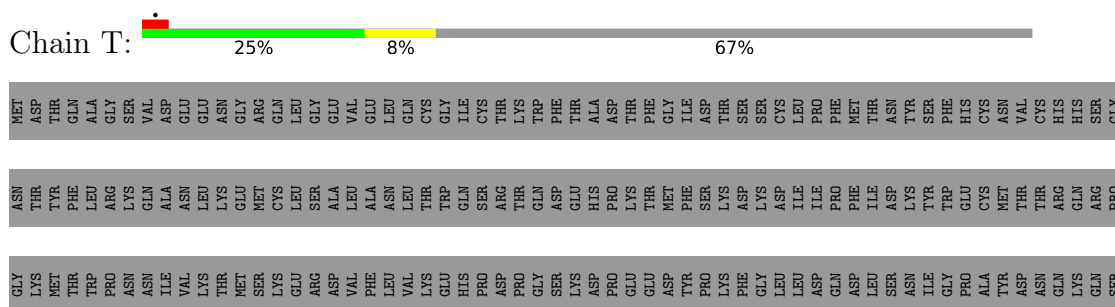
- Molecule 8: Histone-lysine N-methyltransferase 2A

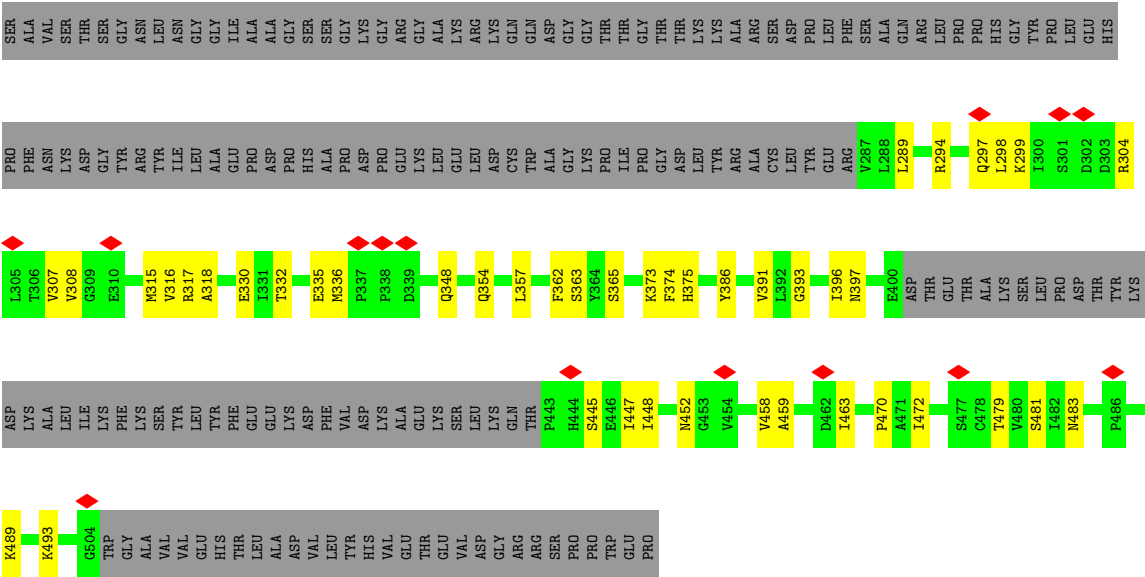


- Molecule 9: WD repeat-containing protein 5



- Molecule 10: Set1/Ash2 histone methyltransferase complex subunit ASH2





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	31882	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	45.0	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.044	Depositor
Minimum map value	-0.019	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.005	Depositor
Map size ($\text{\AA}$)	297.0, 297.0, 297.0	wwPDB
Map dimensions	270, 270, 270	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.37	0/814	0.56	0/1092
1	E	0.37	0/814	0.56	0/1092
2	B	0.39	0/664	0.60	0/889
2	F	0.39	0/629	0.55	0/843
3	C	0.35	0/835	0.52	0/1127
3	G	0.35	0/819	0.53	0/1106
4	D	0.36	0/749	0.52	0/1008
4	H	0.36	0/753	0.52	0/1012
5	I	0.42	0/3311	0.53	0/5103
6	J	0.42	0/3357	0.50	0/5184
7	N	0.23	0/2817	0.53	2/3826 (0.1%)
8	K	0.20	0/1462	0.48	0/1952
9	R	0.18	0/2382	0.43	0/3231
10	T	0.19	0/1442	0.44	0/1950
All	All	0.33	0/20848	0.51	2/29415 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
7	N	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	N	323	GLN	CA-C-N	5.57	134.73	124.82
7	N	323	GLN	C-N-CA	5.57	134.73	124.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
7	N	194	GLY	Peptide

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	802	0	841	8	0
1	E	802	0	841	4	0
2	B	657	0	706	9	0
2	F	622	0	660	2	0
3	C	825	0	882	3	0
3	G	809	0	864	2	0
4	D	738	0	760	3	0
4	H	742	0	765	7	0
5	I	2955	0	1627	13	0
6	J	2990	0	1628	15	0
7	N	2759	0	2707	65	0
8	K	1436	0	1427	32	0
9	R	2326	0	2309	45	0
10	T	1399	0	1363	26	0
11	K	26	0	19	5	0
12	K	1	0	0	0	0
All	All	19889	0	17399	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)
8:K:3907:HIS:H	11:K:4001:SAH:HN61	1.41	0.69
9:R:115:ASP:HB3	9:R:119:GLY:H	1.59	0.68
10:T:297:GLN:HE22	10:T:354:GLN:HE22	1.42	0.68
9:R:111:LEU:HB2	9:R:125:LEU:HB2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:N:12:ASN:HB3	7:N:324:ASN:HD22	1.60	0.66
9:R:41:LEU:HB2	9:R:327:ILE:HB	1.77	0.66
7:N:49:VAL:HG13	7:N:61:ILE:HG12	1.79	0.65
8:K:3861:ASN:HD21	8:K:3897:MET:HB2	1.63	0.64
3:C:15:LYS:HB2	3:C:20:ARG:HH21	1.63	0.62
1:A:42:ARG:NH2	5:I:144:DC:OP2	2.33	0.61
7:N:302:ALA:HB3	7:N:311:ALA:HB3	1.82	0.60
7:N:48:ILE:HB	7:N:62:ILE:HB	1.83	0.60
9:R:69:ILE:HB	9:R:83:ILE:HB	1.83	0.59
10:T:396:ILE:HD11	10:T:463:ILE:HG21	1.84	0.59
7:N:261:GLY:HA3	7:N:305:PRO:HA	1.83	0.59
7:N:27:ALA:H	7:N:316:GLY:HA2	1.68	0.59
7:N:32:PHE:HB3	7:N:36:GLY:HA2	1.85	0.59
4:H:119:THR:HA	4:H:122:LYS:HE3	1.83	0.59
8:K:3906:ASN:ND2	8:K:3939:GLU:OE2	2.36	0.59
5:I:50:DG:N2	6:J:99:DT:O2	2.36	0.59
7:N:280:GLU:HB2	7:N:284:GLY:H	1.68	0.59
8:K:3765:ARG:HD3	9:R:305:ILE:HG21	1.84	0.58
9:R:136:ASN:ND2	9:R:177:VAL:O	2.37	0.58
9:R:110:THR:HG22	9:R:126:LYS:HG2	1.84	0.58
8:K:3849:ILE:HB	8:K:3934:ILE:HB	1.86	0.58
8:K:3864:ARG:NH1	10:T:354:GLN:O	2.36	0.58
9:R:113:ILE:HD12	9:R:123:LYS:HB2	1.86	0.58
8:K:3765:ARG:NH2	9:R:91:SER:O	2.36	0.57
4:D:27:ARG:NH2	5:I:105:DT:OP1	2.37	0.57
9:R:39:PHE:HB2	9:R:329:LEU:HD12	1.86	0.57
9:R:283:VAL:HB	9:R:297:LEU:HB2	1.86	0.57
1:E:108:ASN:ND2	2:F:42:GLY:O	2.38	0.57
7:N:30:CYS:HB2	7:N:39:LEU:HD11	1.86	0.56
9:R:154:ARG:NH1	9:R:166:THR:OG1	2.39	0.56
8:K:3857:GLU:O	8:K:3900:ASN:ND2	2.38	0.56
9:R:179:PHE:HB3	9:R:183:GLY:HA2	1.88	0.56
1:A:108:ASN:ND2	2:B:42:GLY:O	2.39	0.55
10:T:330:GLU:HB2	10:T:483:ASN:HB3	1.88	0.55
1:A:116:ARG:NH1	1:A:118:THR:O	2.40	0.55
8:K:3864:ARG:HA	8:K:3891:GLU:HG2	1.87	0.55
9:R:157:ASP:HB3	9:R:161:GLY:H	1.72	0.55
1:E:116:ARG:NH1	1:E:118:THR:O	2.40	0.55
8:K:3862:VAL:HG22	8:K:3893:VAL:HG22	1.88	0.55
7:N:70:CYS:SG	7:N:71:SER:N	2.80	0.54
7:N:94:ASP:HB3	7:N:98:GLY:H	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:R:65:ALA:HA	9:R:90:ILE:H	1.73	0.54
3:G:71:ARG:NH2	4:H:48:ASP:OD2	2.40	0.54
9:R:278:SER:H	9:R:304:VAL:HB	1.73	0.54
9:R:52:LYS:HD3	9:R:308:ALA:HB2	1.90	0.54
8:K:3778:PHE:HB2	8:K:3841:ARG:HH22	1.72	0.53
9:R:153:VAL:HB	9:R:167:LEU:HB2	1.90	0.53
1:A:106:ASP:OD2	1:A:131:ARG:NH2	2.40	0.53
1:E:106:ASP:OD2	1:E:131:ARG:NH2	2.40	0.53
10:T:315:MET:HE3	10:T:317:ARG:HD3	1.88	0.53
9:R:189:SER:HB3	9:R:220:VAL:HG22	1.91	0.53
1:A:83:ARG:HB2	2:B:80:THR:HG22	1.90	0.53
8:K:3814:LEU:O	8:K:3819:ARG:NH1	2.42	0.53
10:T:336:MET:SD	10:T:386:TYR:OH	2.63	0.53
5:I:8:DA:H2"	5:I:9:DT:H5"	1.91	0.53
2:B:98:TYR:OH	4:H:65:ASP:OD2	2.27	0.52
11:K:4001:SAH:SD	11:K:4001:SAH:N	2.77	0.52
9:R:100:ASN:HA	9:R:116:VAL:HB	1.91	0.52
7:N:33:ASN:OD1	7:N:37:THR:N	2.42	0.52
7:N:259:PHE:HB3	7:N:263:GLY:HA2	1.92	0.52
2:B:23:ARG:HH21	7:N:193:THR:HB	1.74	0.52
7:N:313:ILE:HG12	7:N:318:VAL:HG22	1.91	0.52
9:R:138:ASN:ND2	9:R:142:ASN:OD1	2.40	0.52
8:K:3861:ASN:OD1	8:K:3898:HIS:ND1	2.41	0.51
7:N:67:HIS:NE2	7:N:87:ASP:OD2	2.43	0.51
7:N:83:SER:OG	7:N:93:TRP:NE1	2.44	0.51
10:T:294:ARG:HB2	10:T:316:VAL:HG23	1.91	0.51
7:N:91:SER:HB3	7:N:101:ASP:H	1.76	0.51
9:R:149:PHE:HA	9:R:173:PRO:HB3	1.92	0.51
7:N:76:ARG:NH2	7:N:117:PRO:O	2.42	0.51
7:N:112:LYS:HB3	7:N:126:CYS:HB2	1.92	0.51
10:T:335:GLU:HB3	10:T:479:THR:HB	1.92	0.51
1:E:63:ARG:HH21	6:J:60:DA:H5"	1.76	0.51
10:T:330:GLU:HA	10:T:393:GLY:HA3	1.93	0.51
4:D:33:SER:OG	4:D:34:TYR:N	2.41	0.51
9:R:304:VAL:HA	9:R:320:ALA:HA	1.93	0.51
10:T:332:THR:HB	10:T:481:SER:HB2	1.93	0.51
10:T:289:LEU:O	10:T:304:ARG:NH2	2.38	0.50
7:N:50:ILE:HD12	7:N:60:LYS:HB3	1.92	0.50
9:R:220:VAL:HG12	9:R:231:ALA:HA	1.92	0.50
7:N:235:ARG:NH1	7:N:236:ASP:OD2	2.44	0.50
10:T:447:ILE:HG22	10:T:459:ALA:HB3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:T:308:VAL:HG22	10:T:479:THR:HA	1.93	0.50
2:B:39:ARG:NH1	2:B:43:VAL:O	2.45	0.50
7:N:265:TYR:HA	7:N:280:GLU:HA	1.94	0.50
8:K:3823:LEU:HD13	8:K:3854:MET:HB3	1.94	0.49
10:T:318:ALA:HB3	10:T:470:PRO:HG2	1.94	0.49
1:A:76:GLN:HE22	2:B:21:VAL:HG13	1.78	0.49
2:F:77:LYS:HE3	4:H:89:ARG:HH12	1.77	0.49
7:N:227:GLY:HA2	7:N:230:ILE:HD12	1.94	0.49
9:R:111:LEU:N	9:R:125:LEU:O	2.44	0.49
8:K:3816:MET:HG3	8:K:3920:ILE:HD11	1.94	0.49
8:K:3859:ALA:H	8:K:3900:ASN:HD21	1.60	0.49
6:J:94:DG:H2'	6:J:95:DG:C8	2.47	0.49
10:T:307:VAL:HG11	10:T:472:ILE:HD13	1.94	0.49
7:N:113:VAL:HG12	7:N:125:VAL:HG22	1.93	0.49
7:N:255:LYS:HB2	7:N:271:ALA:HB2	1.95	0.49
1:A:61:LEU:HD11	2:B:40:ARG:HH21	1.77	0.48
7:N:88:ASN:HB3	7:N:106:PHE:HD2	1.77	0.48
9:R:238:LEU:HB2	9:R:252:TYR:HD2	1.78	0.48
6:J:35:DT:H2''	6:J:36:DA:C8	2.47	0.48
7:N:260:SER:HB3	7:N:265:TYR:HB2	1.94	0.48
7:N:340:ASP:O	8:K:3824:LYS:NZ	2.43	0.48
5:I:60:DA:N6	6:J:87:DT:O4	2.46	0.48
9:R:195:CYS:HB2	9:R:209:LEU:HB3	1.95	0.48
3:C:68:ASN:OD1	3:C:71:ARG:NH1	2.46	0.48
5:I:92:DC:C2	6:J:56:DG:N2	2.81	0.48
6:J:94:DG:H2'	6:J:95:DG:H8	1.79	0.48
7:N:75:SER:OG	7:N:77:ASP:OD1	2.31	0.48
9:R:88:LEU:HD12	9:R:108:ASP:HA	1.96	0.48
9:R:155:ILE:HG22	9:R:164:LEU:HD12	1.96	0.48
7:N:50:ILE:N	7:N:60:LYS:O	2.43	0.47
8:K:3863:ILE:HD13	8:K:3871:ARG:HD2	1.96	0.47
2:B:47:SER:OG	2:B:48:GLY:N	2.46	0.47
7:N:223:ARG:HG2	7:N:244:LYS:HG2	1.96	0.47
7:N:269:GLY:HA2	7:N:276:LEU:HD22	1.96	0.47
9:R:91:SER:OG	9:R:107:ASP:OD1	2.28	0.47
9:R:90:ILE:HA	9:R:106:SER:HA	1.95	0.47
10:T:365:SER:OG	10:T:374:PHE:N	2.38	0.47
9:R:123:LYS:HE2	9:R:160:THR:HA	1.96	0.47
10:T:357:LEU:HD12	10:T:363:SER:HB2	1.95	0.47
7:N:67:HIS:HB2	7:N:85:SER:HB2	1.97	0.47
10:T:298:LEU:HD12	10:T:316:VAL:HB	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:N:20:THR:OG1	7:N:318:VAL:O	2.30	0.46
9:R:103:VAL:HG23	9:R:113:ILE:HG12	1.97	0.46
7:N:34:ARG:NH1	7:N:76:ARG:O	2.47	0.46
5:I:100:DG:H2''	5:I:101:DG:C8	2.51	0.46
7:N:47:ARG:HG2	7:N:63:SER:HA	1.98	0.46
9:R:275:VAL:HG22	9:R:285:ILE:HG12	1.98	0.46
10:T:373:LYS:NZ	10:T:459:ALA:O	2.48	0.46
9:R:186:ILE:HB	9:R:198:TRP:HB2	1.97	0.46
9:R:286:TRP:CD1	9:R:294:VAL:H	2.33	0.46
10:T:397:ASN:HB3	10:T:445:SER:HA	1.98	0.46
5:I:66:DC:H2''	5:I:67:DG:C8	2.51	0.46
7:N:193:THR:OG1	7:N:223:ARG:NH2	2.49	0.46
7:N:300:ASP:HB3	7:N:313:ILE:HB	1.98	0.46
3:C:14:ALA:HB1	5:I:31:DT:H4'	1.98	0.45
4:D:115:VAL:O	4:D:119:THR:OG1	2.33	0.45
3:G:42:ARG:HH11	5:I:112:DT:H4'	1.81	0.45
6:J:115:DC:N3	6:J:116:DA:N6	2.65	0.45
8:K:3861:ASN:HB3	8:K:3863:ILE:HG23	1.98	0.45
4:H:83:ARG:HH21	6:J:40:DA:H3'	1.81	0.45
7:N:276:LEU:N	7:N:290:LEU:O	2.37	0.45
8:K:3861:ASN:N	8:K:3894:ASP:O	2.49	0.45
7:N:91:SER:OG	7:N:102:GLN:O	2.25	0.45
7:N:202:LYS:NZ	7:N:254:TRP:O	2.39	0.45
6:J:57:DT:H2''	6:J:58:DT:C5	2.51	0.45
10:T:452:ASN:O	10:T:493:LYS:NZ	2.50	0.45
9:R:45:THR:OG1	9:R:66:ASP:OD2	2.35	0.45
7:N:185:LEU:HD11	7:N:188:SER:HB3	1.99	0.45
7:N:278:ILE:HD12	7:N:288:LYS:HB3	1.98	0.44
10:T:299:LYS:HB2	10:T:308:VAL:HB	1.98	0.44
9:R:57:GLY:O	9:R:310:HIS:NE2	2.45	0.44
9:R:229:ILE:O	9:R:241:TRP:N	2.41	0.44
7:N:74:TRP:HA	7:N:81:LEU:HD23	1.99	0.44
7:N:123:VAL:O	7:N:135:LEU:N	2.41	0.44
8:K:3883:TYR:OH	11:K:4001:SAH:SD	2.66	0.44
5:I:24:DC:H2''	5:I:25:DG:C8	2.53	0.44
7:N:22:ASP:OD1	7:N:318:VAL:N	2.51	0.44
7:N:94:ASP:H	7:N:98:GLY:HA2	1.82	0.44
6:J:82:DG:H2'	6:J:83:DT:C6	2.53	0.44
7:N:134:MET:O	7:N:143:VAL:N	2.41	0.43
8:K:3836:SER:OG	8:K:3906:ASN:ND2	2.51	0.43
8:K:3838:ILE:O	8:K:3958:ASN:ND2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:R:199:ASP:HB3	9:R:204:GLN:H	1.82	0.43
6:J:78:DG:H4'	6:J:79:DT:H5'	1.99	0.43
7:N:121:ASN:O	7:N:137:LEU:N	2.51	0.43
7:N:217:THR:OG1	7:N:219:ASP:OD1	2.26	0.43
8:K:3835:ARG:HA	8:K:3841:ARG:HG3	2.00	0.43
10:T:348:GLN:HB2	10:T:362:PHE:HB3	2.00	0.43
7:N:260:SER:N	7:N:265:TYR:O	2.51	0.43
8:K:3846:LYS:HA	8:K:3936:ARG:HG3	2.01	0.43
7:N:51:TRP:HA	7:N:59:ALA:H	1.83	0.43
8:K:3810:THR:HB	8:K:3854:MET:H	1.84	0.43
4:H:48:ASP:OD1	4:H:48:ASP:N	2.50	0.43
7:N:14:PRO:HB3	7:N:308:PRO:HB2	1.99	0.43
8:K:3840:GLY:HA3	11:K:4001:SAH:HB2	2.01	0.43
8:K:3909:CYS:H	8:K:3959:CYS:HB2	1.82	0.43
7:N:15:GLU:H	7:N:323:GLN:HB2	1.83	0.43
10:T:363:SER:O	10:T:375:HIS:ND1	2.45	0.43
5:I:82:DC:H2''	5:I:83:DG:C8	2.53	0.42
7:N:299:LEU:HD12	7:N:315:SER:H	1.83	0.42
9:R:301:THR:OG1	9:R:323:ASN:O	2.34	0.42
1:A:69:ARG:NH1	7:N:196:SER:OG	2.52	0.42
5:I:68:DT:H2''	5:I:69:DA:N7	2.34	0.42
7:N:133:VAL:HG22	7:N:144:VAL:HG22	2.02	0.42
6:J:9:DT:H2''	6:J:10:DA:C8	2.55	0.42
9:R:262:ILE:HD11	9:R:279:GLU:HG3	2.02	0.42
2:B:64:ASN:OD1	2:B:67:ARG:NH2	2.52	0.41
7:N:49:VAL:HG22	7:N:61:ILE:HG23	2.02	0.41
8:K:3865:SER:O	8:K:3868:THR:OG1	2.33	0.41
6:J:95:DG:H2'	6:J:96:DT:H71	2.03	0.41
11:K:4001:SAH:H8	11:K:4001:SAH:H2'	1.88	0.41
8:K:3861:ASN:O	8:K:3894:ASP:N	2.43	0.41
8:K:3861:ASN:HB2	8:K:3894:ASP:HB3	2.02	0.41
9:R:240:LEU:HD22	9:R:288:LEU:HD11	2.03	0.41
10:T:448:ILE:HG12	10:T:458:VAL:HA	2.02	0.41
7:N:276:LEU:O	7:N:290:LEU:N	2.53	0.41
10:T:391:VAL:HG21	10:T:489:LYS:HG2	2.02	0.41
7:N:137:LEU:HD23	7:N:137:LEU:HA	1.95	0.41
7:N:208:ARG:HB2	7:N:262:ASP:HA	2.03	0.41
7:N:377:VAL:HG13	7:N:378:THR:HG23	2.02	0.41
9:R:274:ILE:N	9:R:286:TRP:O	2.50	0.41
9:R:321:LEU:HB3	9:R:322:GLU:H	1.74	0.41
4:H:77:LEU:HD23	4:H:77:LEU:HA	1.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:N:88:ASN:ND2	7:N:106:PHE:O	2.53	0.41
6:J:5:DG:H2"	6:J:6:DA:C8	2.56	0.40
8:K:3816:MET:HE2	8:K:3819:ARG:HH21	1.87	0.40
7:N:14:PRO:HG2	7:N:310:ILE:HD11	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	95/135 (70%)	93 (98%)	2 (2%)	0	100	100
1	E	95/135 (70%)	93 (98%)	2 (2%)	0	100	100
2	B	80/102 (78%)	76 (95%)	4 (5%)	0	100	100
2	F	76/102 (74%)	71 (93%)	5 (7%)	0	100	100
3	C	105/129 (81%)	101 (96%)	4 (4%)	0	100	100
3	G	103/129 (80%)	100 (97%)	3 (3%)	0	100	100
4	D	92/122 (75%)	89 (97%)	3 (3%)	0	100	100
4	H	93/122 (76%)	91 (98%)	2 (2%)	0	100	100
7	N	347/538 (64%)	314 (90%)	33 (10%)	0	100	100
8	K	173/216 (80%)	156 (90%)	17 (10%)	0	100	100
9	R	298/334 (89%)	274 (92%)	24 (8%)	0	100	100
10	T	172/534 (32%)	155 (90%)	17 (10%)	0	100	100
All	All	1729/2598 (67%)	1613 (93%)	116 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	85/110 (77%)	85 (100%)	0	100	100
1	E	85/110 (77%)	85 (100%)	0	100	100
2	B	68/78 (87%)	68 (100%)	0	100	100
2	F	64/78 (82%)	64 (100%)	0	100	100
3	C	84/101 (83%)	84 (100%)	0	100	100
3	G	83/101 (82%)	83 (100%)	0	100	100
4	D	81/102 (79%)	81 (100%)	0	100	100
4	H	81/102 (79%)	81 (100%)	0	100	100
7	N	308/462 (67%)	308 (100%)	0	100	100
8	K	151/189 (80%)	151 (100%)	0	100	100
9	R	262/291 (90%)	262 (100%)	0	100	100
10	T	149/460 (32%)	149 (100%)	0	100	100
All	All	1501/2184 (69%)	1501 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	76	GLN
3	C	38	ASN
4	D	46	HIS
1	E	39	HIS
3	G	89	ASN
3	G	110	ASN
4	H	46	HIS
7	N	12	ASN
7	N	154	ASN
7	N	243	GLN

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Mol	Chain	Res	Type
7	N	323	GLN
7	N	324	ASN
7	N	328	ASN
8	K	3906	ASN
8	K	3907	HIS
10	T	292	HIS
10	T	354	GLN
10	T	381	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
11	SAH	K	4001	-	27,28,28	1.10	4 (14%)	36,40,40	2.13	10 (27%)

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
11	SAH	K	4001	-	-	10/15/31/31	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	K	4001	SAH	C2-N3	2.50	1.38	1.33
11	K	4001	SAH	C2-N1	2.41	1.38	1.33
11	K	4001	SAH	OXT-C	-2.35	1.23	1.30
11	K	4001	SAH	C5-N7	-2.15	1.35	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	K	4001	SAH	N3-C2-N1	-5.60	120.10	128.58
11	K	4001	SAH	C5-C4-N3	-4.60	120.39	126.72
11	K	4001	SAH	C5'-SD-CG	-4.31	89.47	102.26
11	K	4001	SAH	N9-C8-N7	-3.94	108.35	113.94
11	K	4001	SAH	C5-N7-C8	3.54	109.01	103.45
11	K	4001	SAH	C2-N3-C4	3.40	120.14	111.83
11	K	4001	SAH	N3-C4-N9	3.19	132.59	127.17
11	K	4001	SAH	OXT-C-O	-3.03	117.20	124.08
11	K	4001	SAH	C4-C5-N7	-2.30	107.95	110.58
11	K	4001	SAH	C3'-C2'-C1'	2.23	105.69	101.46

There are no chirality outliers.

All (10) torsion outliers are listed below:

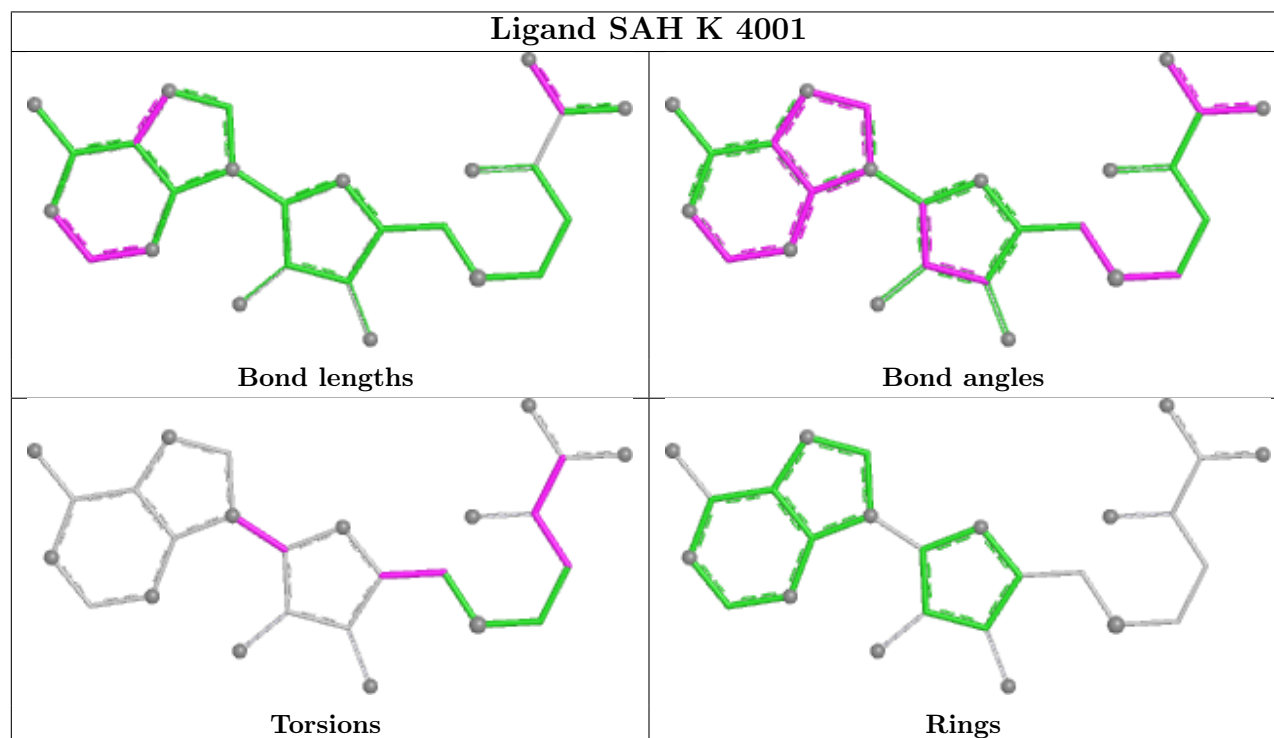
Mol	Chain	Res	Type	Atoms
11	K	4001	SAH	O-C-CA-N
11	K	4001	SAH	O4'-C4'-C5'-SD
11	K	4001	SAH	C3'-C4'-C5'-SD
11	K	4001	SAH	OXT-C-CA-N
11	K	4001	SAH	C2'-C1'-N9-C8
11	K	4001	SAH	C-CA-CB-CG
11	K	4001	SAH	OXT-C-CA-CB
11	K	4001	SAH	O-C-CA-CB
11	K	4001	SAH	C2'-C1'-N9-C4
11	K	4001	SAH	O4'-C1'-N9-C8

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	K	4001	SAH	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

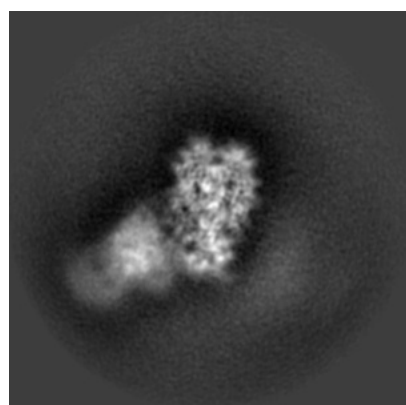
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0695. These allow visual inspection of the internal detail of the map and identification of artifacts.

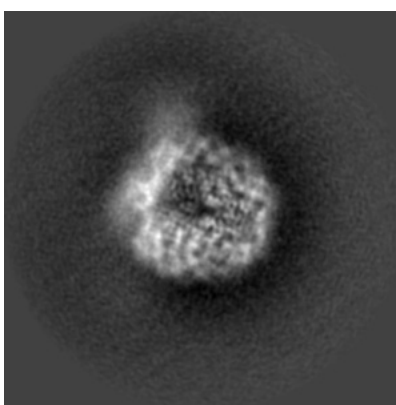
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

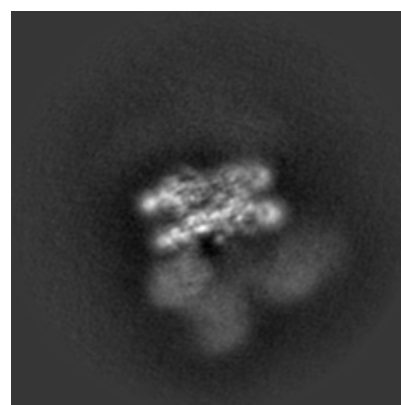
6.1.1 Primary map



X



Y

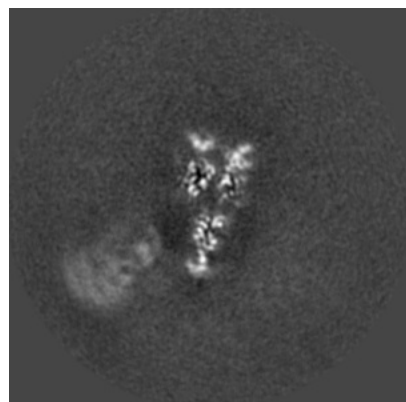


Z

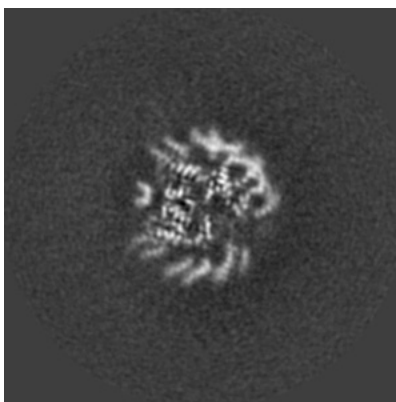
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

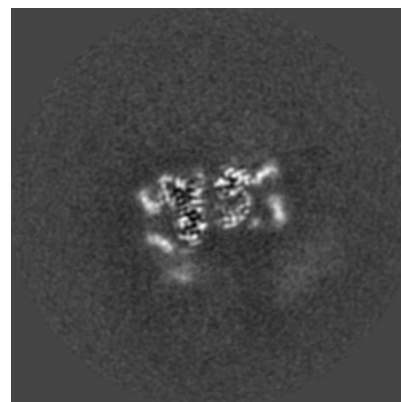
6.2.1 Primary map



X Index: 135



Y Index: 135

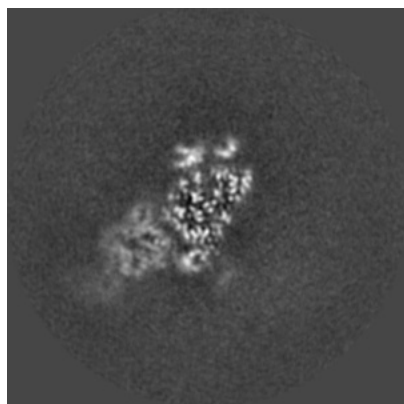


Z Index: 135

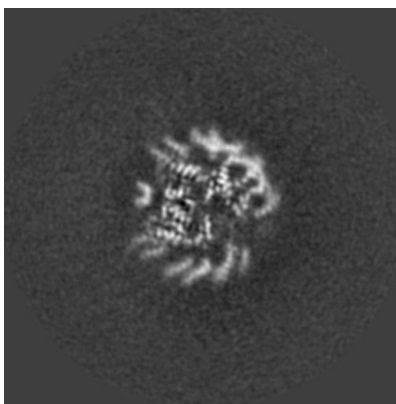
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

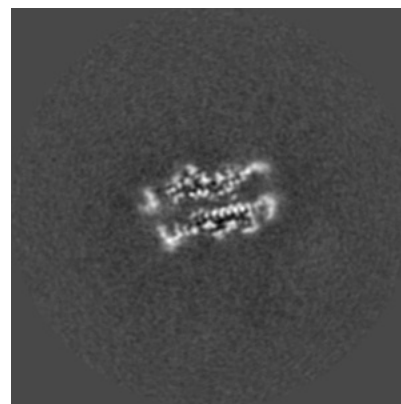
6.3.1 Primary map



X Index: 120



Y Index: 135

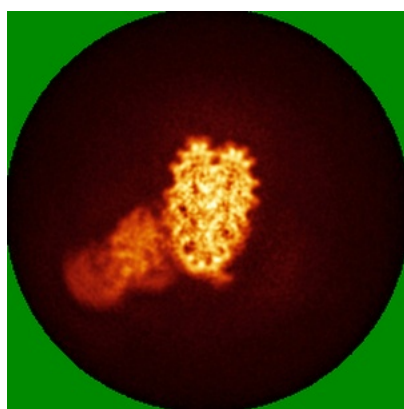


Z Index: 149

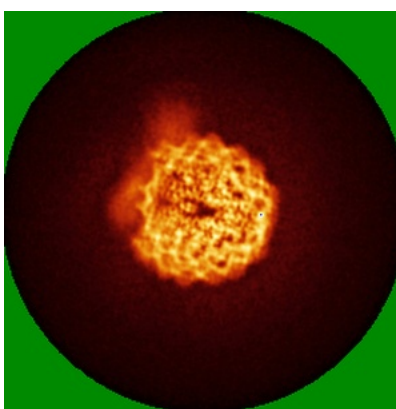
The images above show the largest variance slices of the map in three orthogonal directions.

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

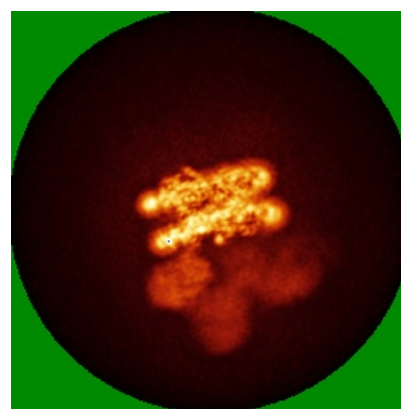
6.4.1 Primary map



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



The images above show the 3D surface view of the map at the recommended contour level 0.005. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

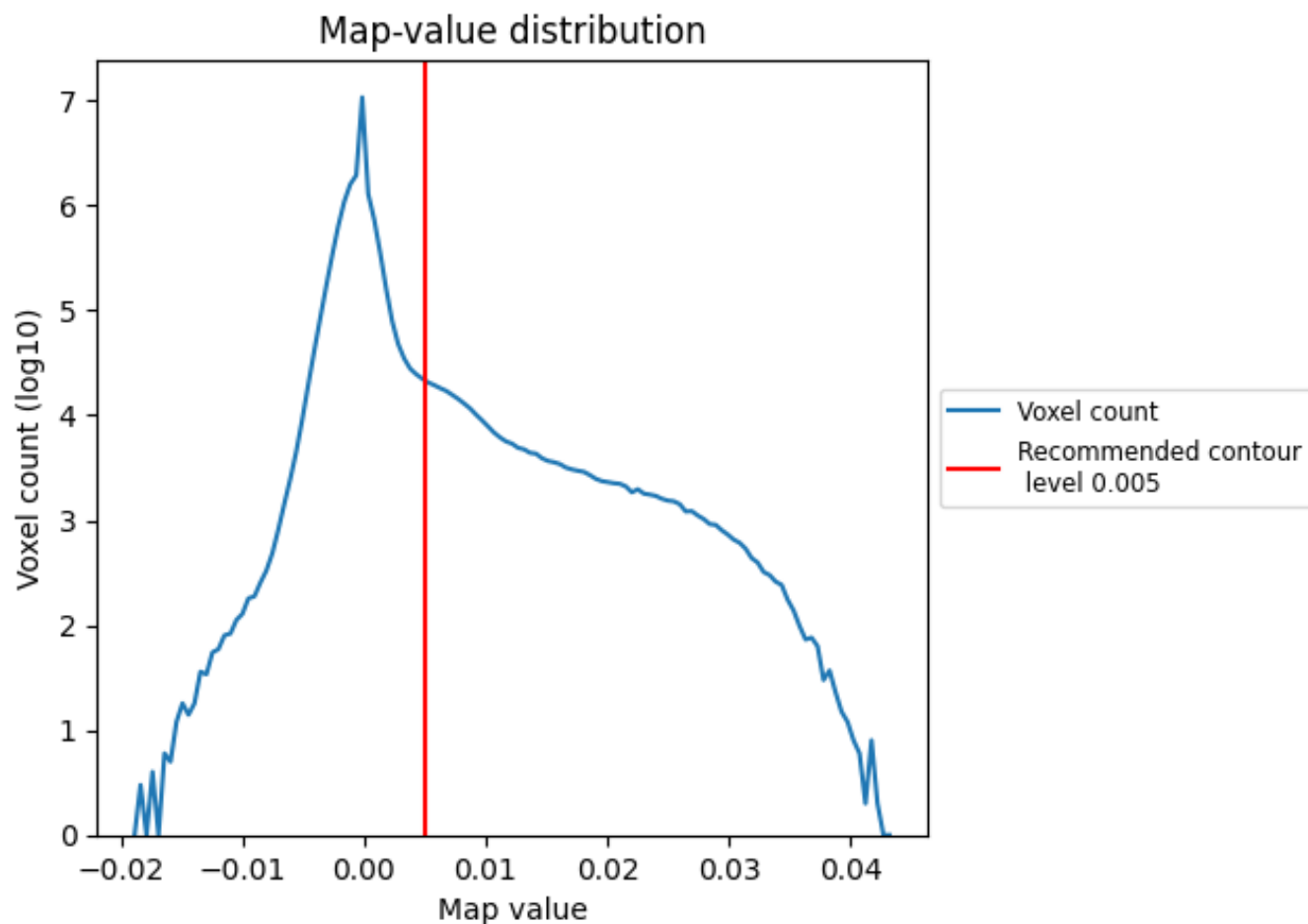
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

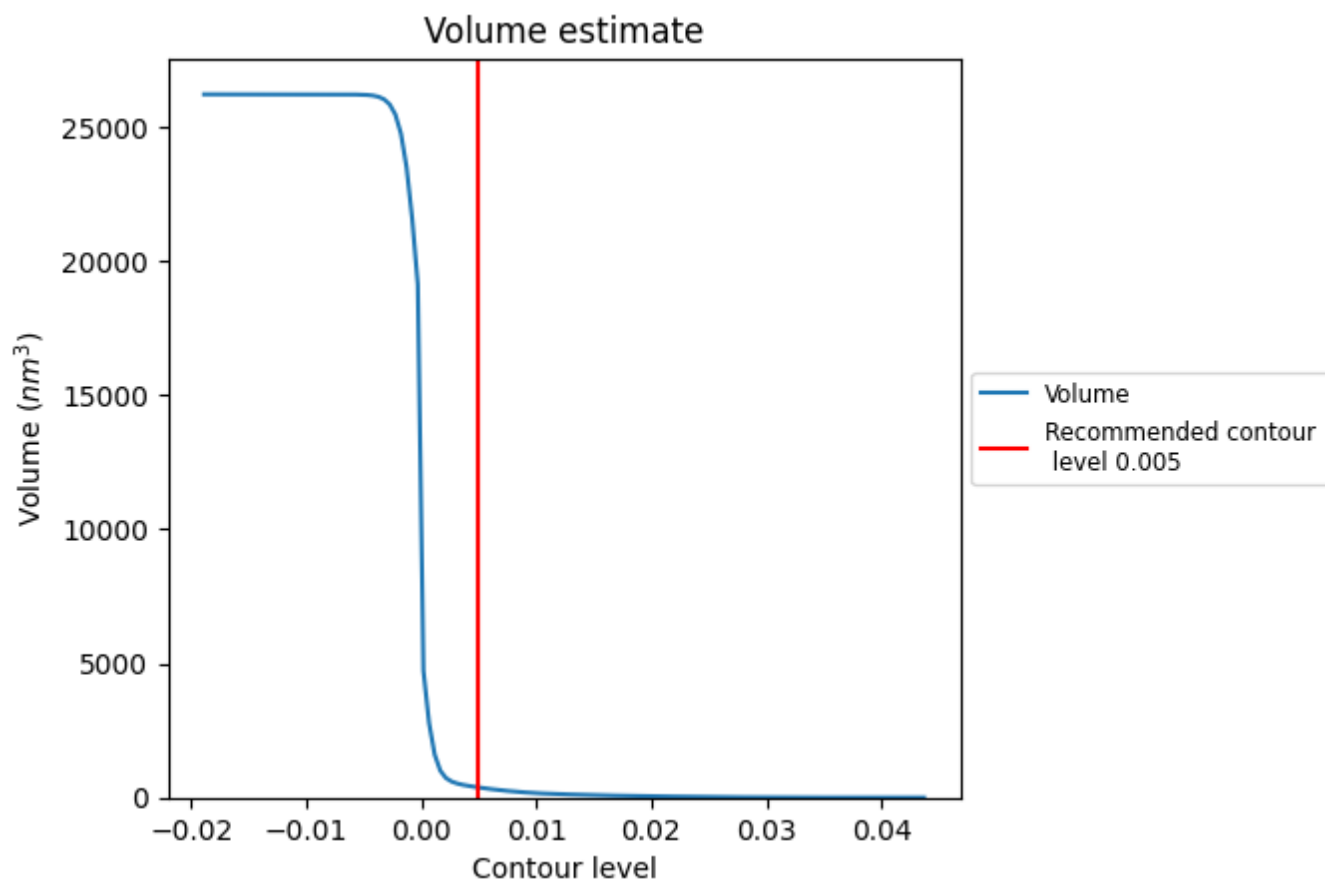
This section contains the results of statistical analysis of the map.

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.

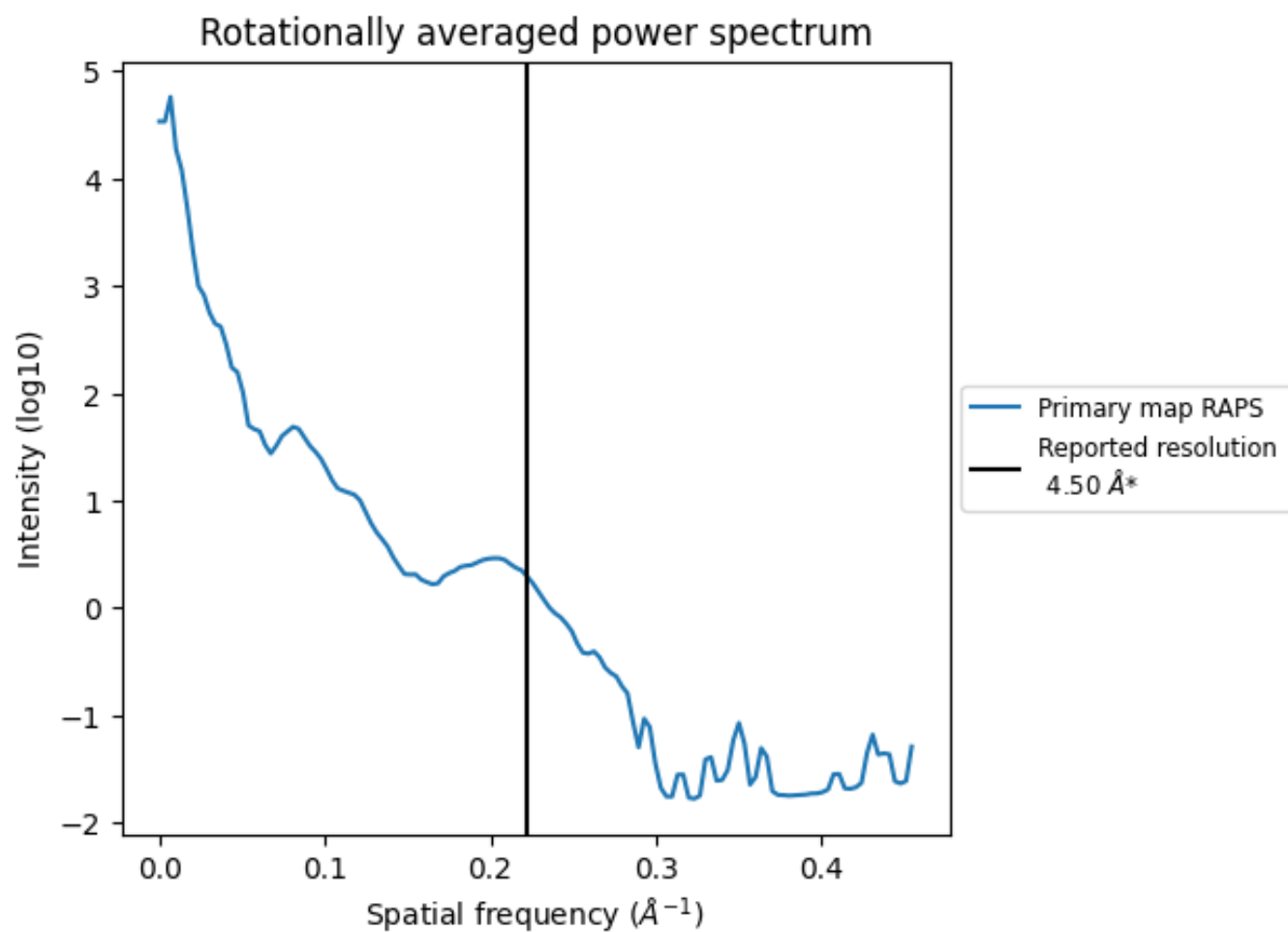
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 375 nm³; this corresponds to an approximate mass of 339 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.222 Å⁻¹

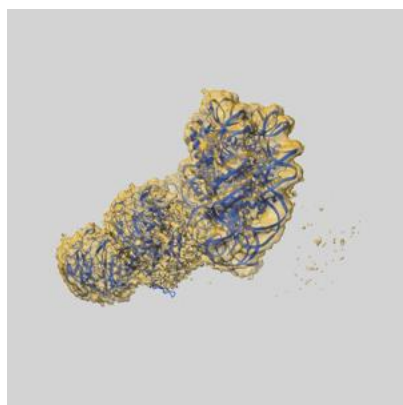
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

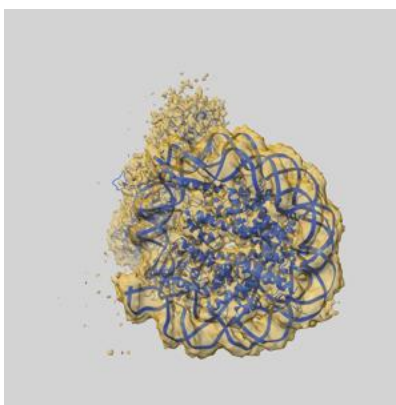
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-0695 and PDB model 6KIZ. Per-residue inclusion information can be found in section [3](#) on page [7](#).

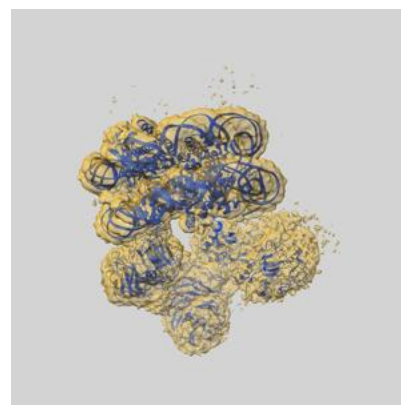
9.1 Map-model overlay [i](#)



X



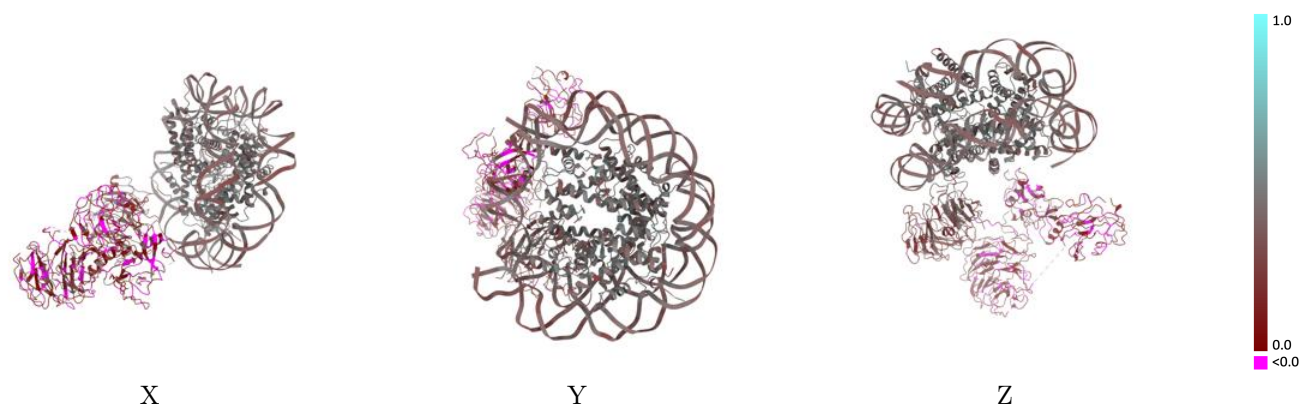
Y



Z

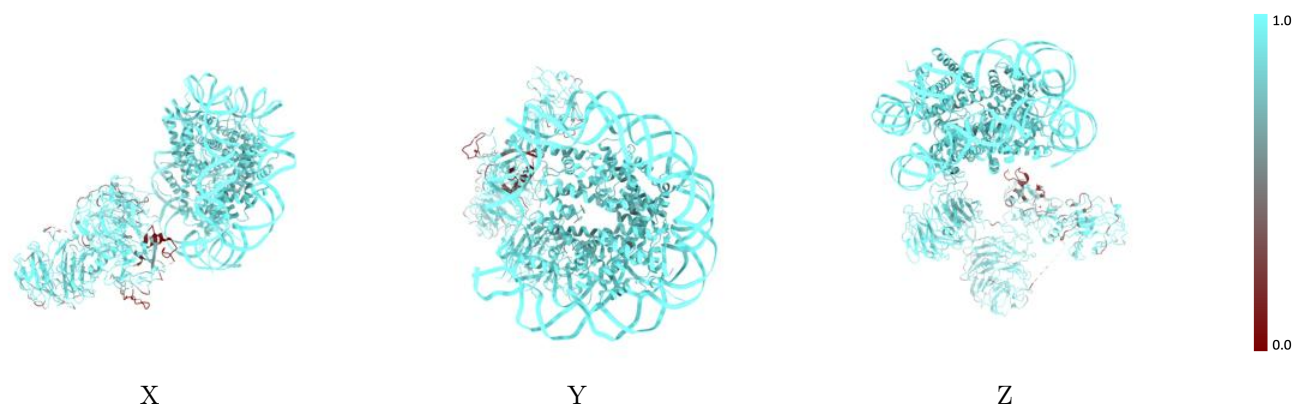
The images above show the 3D surface view of the map at the recommended contour level 0.005 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



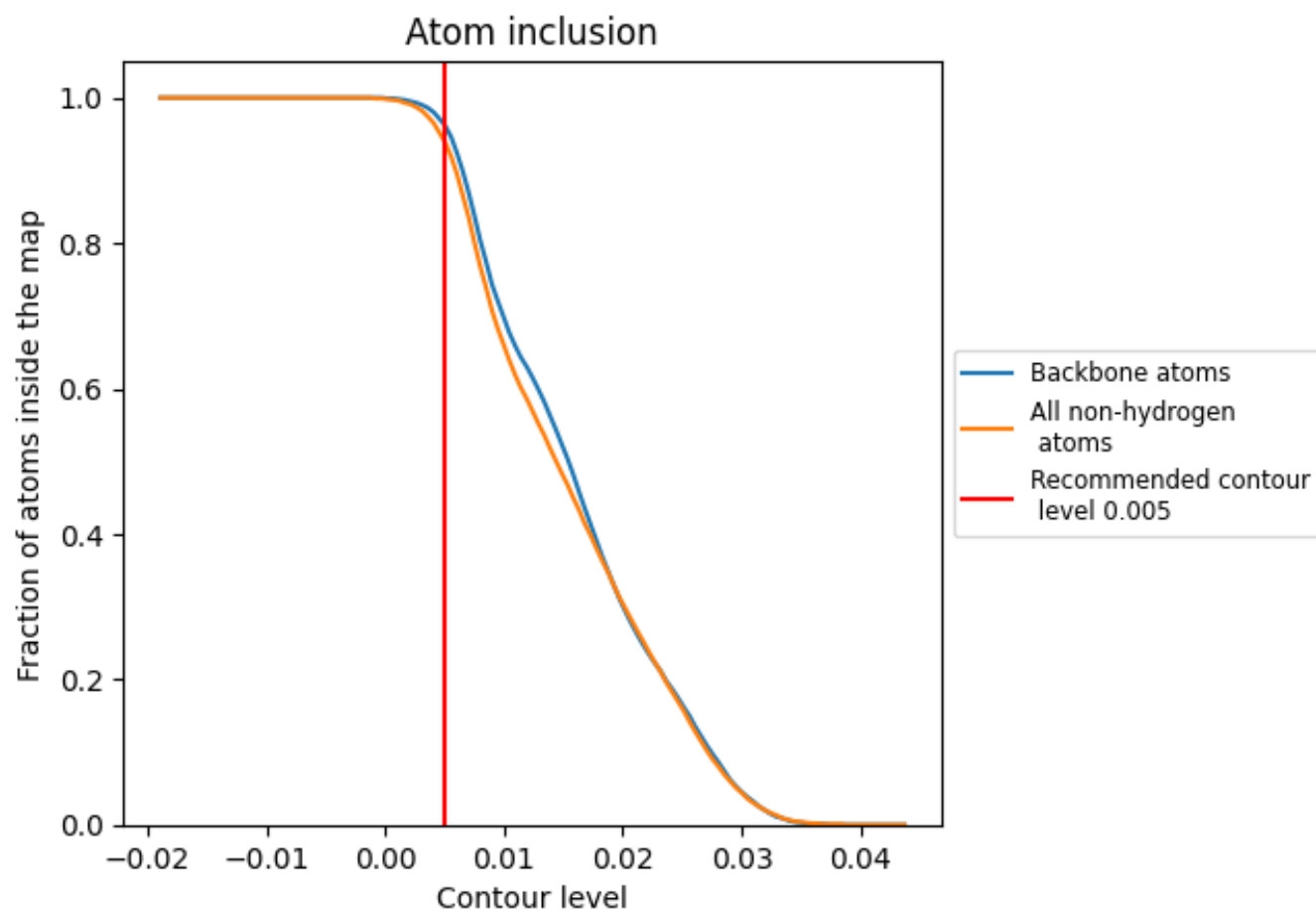
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.005).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 94% 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 (0.005) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9400	 0.3100
A	 0.9510	 0.4450
B	 0.9570	 0.4510
C	 0.9640	 0.4530
D	 0.9650	 0.4310
E	 0.9550	 0.4400
F	 0.9620	 0.4620
G	 0.9710	 0.4460
H	 0.9700	 0.4330
I	 1.0000	 0.3770
J	 1.0000	 0.3790
K	 0.7210	 0.1170
N	 0.9440	 0.2460
R	 0.9100	 0.1160
T	 0.8610	 0.0960

