



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 09:35 AM UTC

PDB ID : 7C79 / pdb_00007c79
EMDB ID : EMD-30296
Title : Cryo-EM structure of yeast Ribonuclease MRP
Authors : Lan, P.; Wu, J.; Lei, M.
Deposited on : 2020-05-24
Resolution : 2.50 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
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

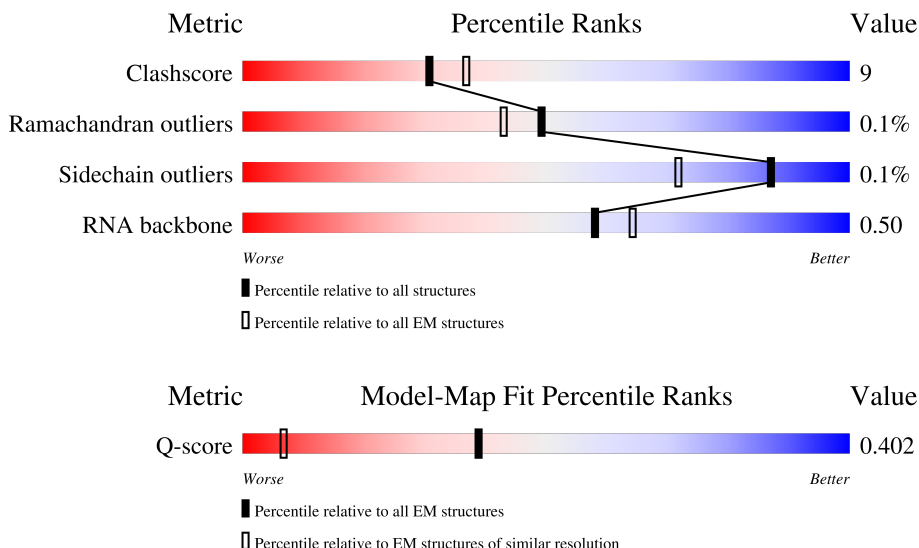
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.50 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	-
RNA backbone	8273	3508	-
Q-score	-	25397	7115 (2.00 - 3.00)

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	340	5% 42% 46% 10% .
2	B	875	. 71% 19% 9%
3	C	195	18% 43% 34% 22%

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Mol	Chain	Length	Quality of chain
4	D	279	 5% 65% 25% 10%
5	E	173	 71% 13% 16%
6	F	158	 84% 15%
7	G	140	 78% 9% 14%
8	H	133	 5% 82% 17%
9	I	293	 63% 19% 17%
9	J	293	 77% 23%
10	K	198	 36% 15% 47%
11	L	201	 50% 15% 35%

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 27324 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 RNA chain called Ribonuclease MRP RNA subunit NME1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	331	Total	C	N	O	P	0	0
			7008	3139	1204	2334	331		

- Molecule 2 is a protein called Ribonucleases P/MRP protein subunit POP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	793	Total	C	N	O	S	0	0
			6453	4085	1170	1162	36		

- Molecule 3 is a protein called Ribonucleases P/MRP protein subunit POP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	152	Total	C	N	O	S	0	0
			1246	812	206	219	9		

- Molecule 4 is a protein called RNases MRP/P 32.9 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	251	Total	C	N	O	S	0	0
			2086	1338	352	386	10		

- Molecule 5 is a protein called Ribonuclease P/MRP protein subunit POP5.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	145	Total	C	N	O	S	0	0
			1135	716	201	210	8		

- Molecule 6 is a protein called Ribonucleases P/MRP protein subunit POP6.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	157	Total	C	N	O	S	0	0
			1271	804	222	241	4		

- Molecule 7 is a protein called Ribonucleases P/MRP protein subunit POP7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	121	Total	C	N	O	S	0	0
			960	609	167	182	2		

- Molecule 8 is a protein called Ribonucleases P/MRP protein subunit POP8.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	131	Total	C	N	O	S	0	0
			1079	683	177	215	4		

- Molecule 9 is a protein called Ribonuclease P/MRP protein subunit RPP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	242	Total	C	N	O	S	0	0
			1881	1188	339	345	9		
9	J	292	Total	C	N	O	S	0	0
			2253	1410	412	421	10		

- Molecule 10 is a protein called Ribonuclease MRP protein subunit SNM1.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	K	104	Total	C	N	O	S	0	0
			840	527	148	156	9		

- Molecule 11 is a protein called Ribonuclease MRP protein subunit RMP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	L	131	Total	C	N	O	S	0	0
			1110	711	210	182	7		

- Molecule 12 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
12	A	1	Total	Mg	0
			1	1	

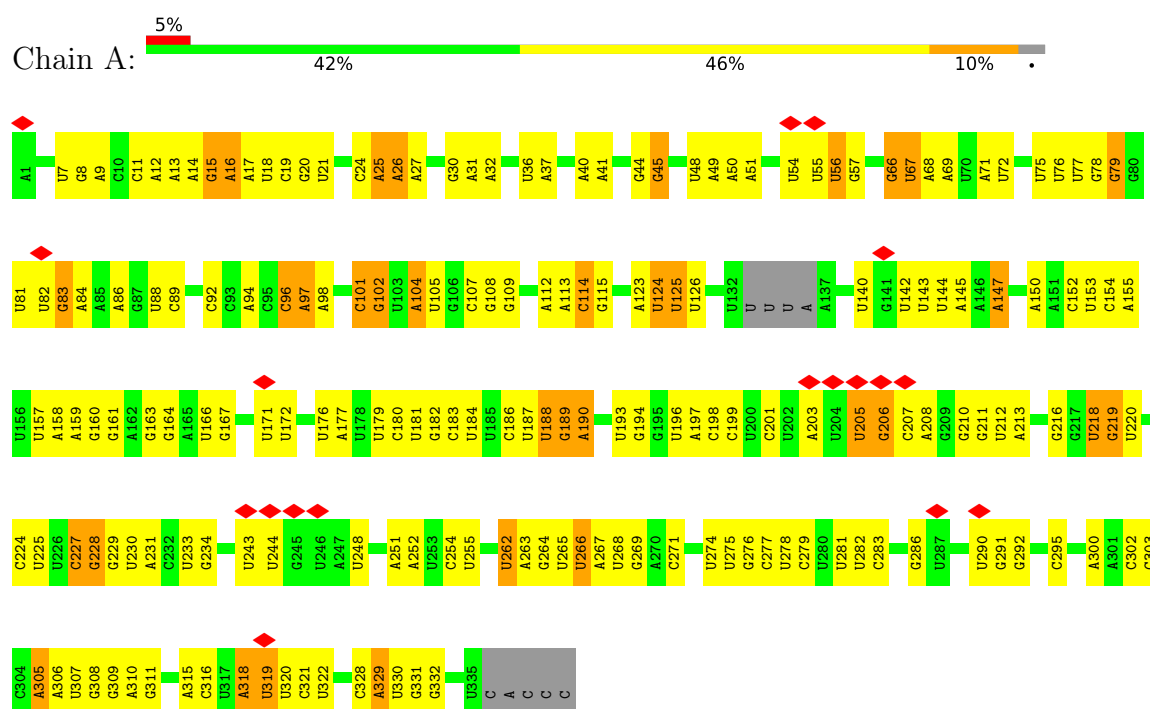
- Molecule 13 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

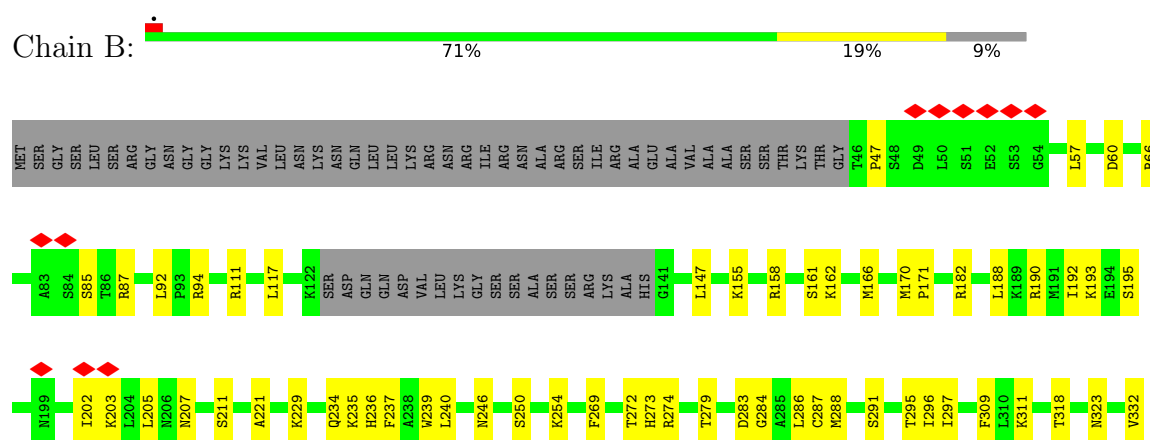
3 Residue-property plots

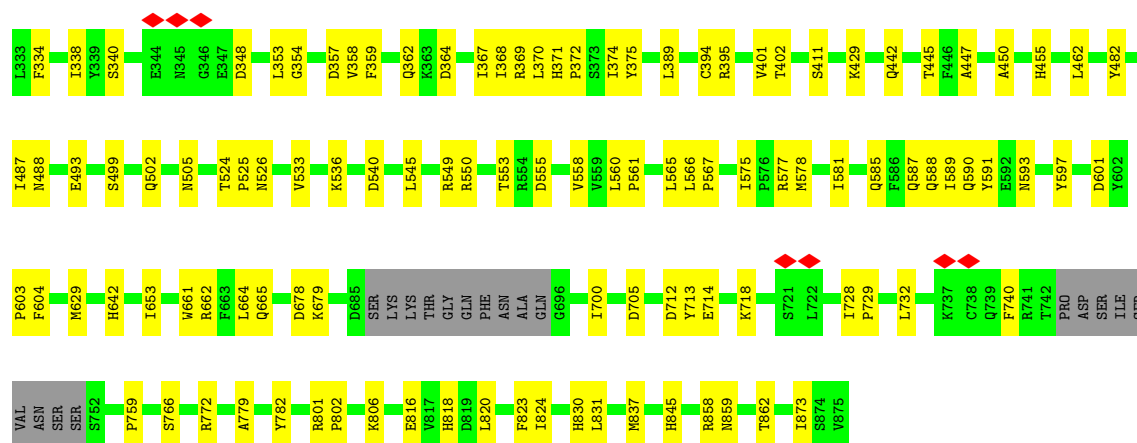
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: Ribonuclease MRP RNA subunit NME1

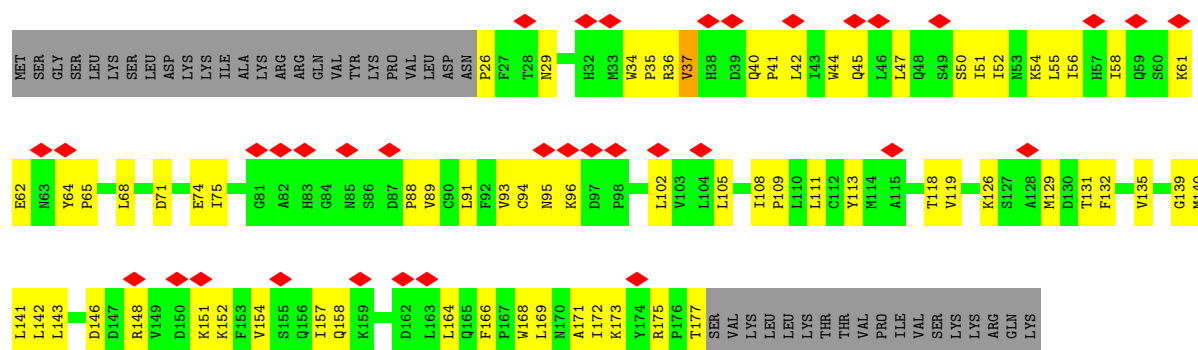
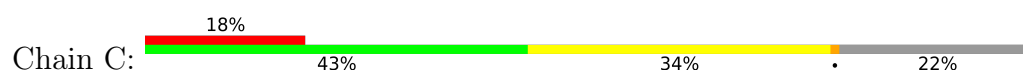


• Molecule 2: Ribonucleases P/MRP protein subunit POP1

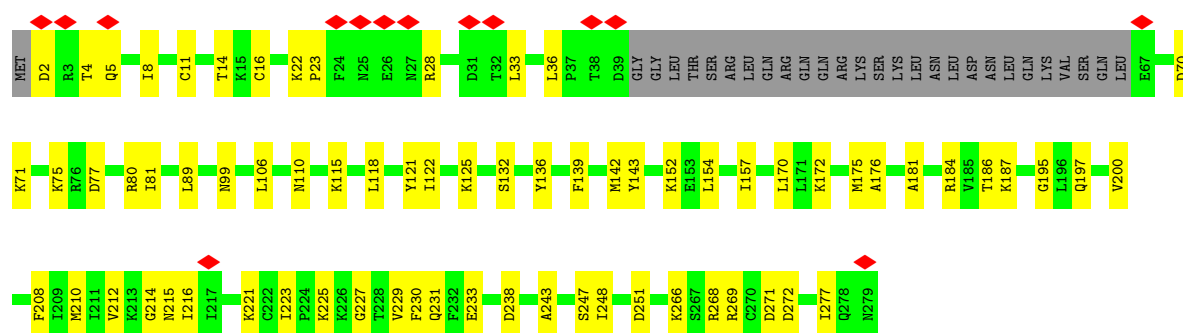




• Molecule 3: Ribonucleases P/MRP protein subunit POP3



• Molecule 4: RNases MRP/P 32.9 kDa subunit



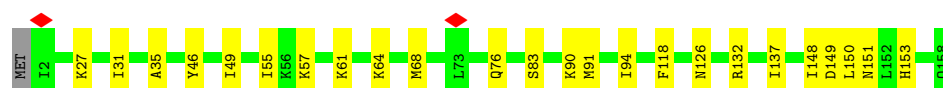
• Molecule 5: Ribonuclease P/MRP protein subunit POP5



ILE
GLY
ARG
GLU
ASN
GLU
ASN
GLU
ASP
ASP

• Molecule 6: Ribonucleases P/MRP protein subunit POP6

Chain F:  84% 15%



• Molecule 7: Ribonucleases P/MRP protein subunit POP7

Chain G:  78% 9% 14%



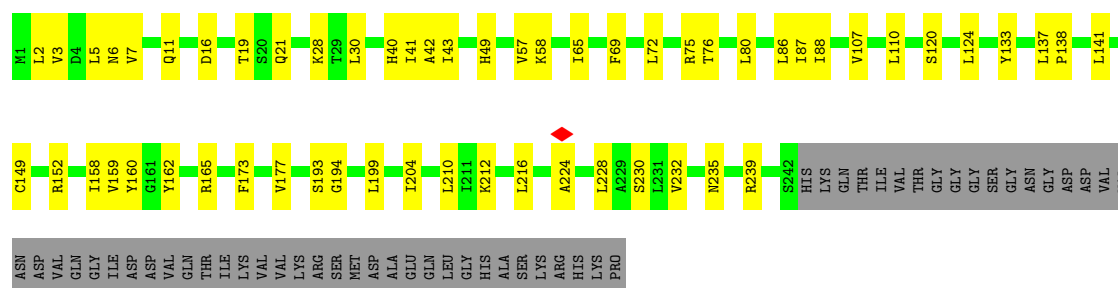
• Molecule 8: Ribonucleases P/MRP protein subunit POP8

Chain H:  5% 82% 17%



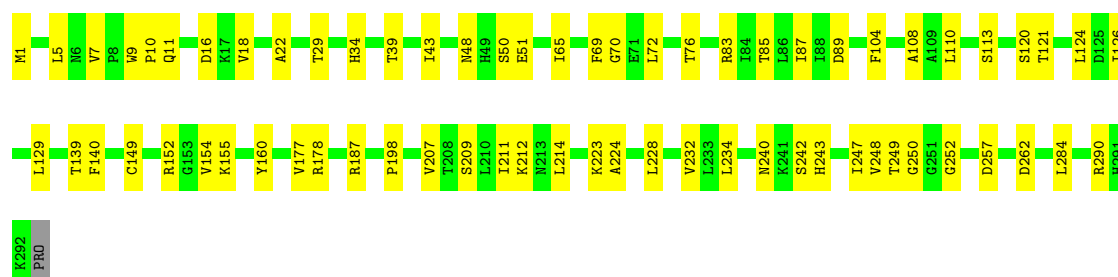
• Molecule 9: Ribonuclease P/MRP protein subunit RPP1

Chain I:  63% 19% 17%

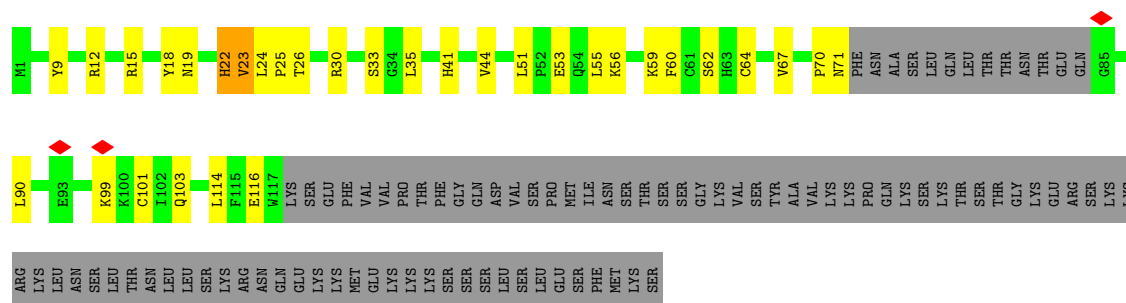
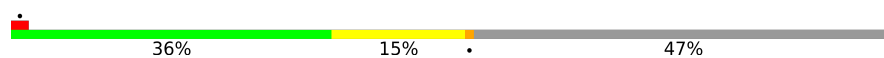


• Molecule 9: Ribonuclease P/MRP protein subunit RPP1

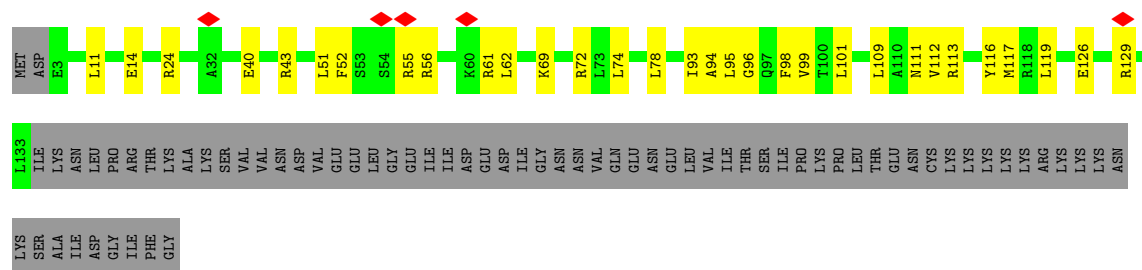
Chain J:  77% 23%



Chain K:



Chain L:



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	1232761	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	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	0.105	Depositor
Minimum map value	-0.042	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.007	Depositor
Map size ($\text{\AA}$)	422.40002, 422.40002, 422.40002	wwPDB
Map dimensions	384, 384, 384	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, MG

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	0.07	0/7828	0.21	0/12182
2	B	0.15	0/6593	0.43	0/8887
3	C	0.25	0/1281	0.64	0/1744
4	D	0.14	0/2123	0.39	0/2851
5	E	0.12	0/1148	0.32	0/1542
6	F	0.10	0/1289	0.31	0/1733
7	G	0.10	0/972	0.32	0/1305
8	H	0.14	0/1101	0.39	0/1488
9	I	0.13	0/1908	0.42	0/2577
9	J	0.12	0/2283	0.37	0/3080
10	K	0.15	0/856	0.53	0/1145
11	L	0.14	0/1132	0.44	0/1523
All	All	0.13	0/28514	0.37	0/40057

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
10	K	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
10	K	22	HIS	Peptide

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Mol	Chain	Res	Type	Group
10	K	99	LYS	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	7008	0	3534	113	0
2	B	6453	0	6530	116	0
3	C	1246	0	1261	52	0
4	D	2086	0	2132	53	0
5	E	1135	0	1199	17	0
6	F	1271	0	1305	16	0
7	G	960	0	1005	7	0
8	H	1079	0	1037	12	0
9	I	1881	0	1975	35	0
9	J	2253	0	2344	44	0
10	K	840	0	830	25	0
11	L	1110	0	1138	22	0
12	A	1	0	0	0	0
13	K	1	0	0	0	0
All	All	27324	0	24290	455	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 9.

All (455) 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)
3:C:151:LYS:HA	3:C:154:VAL:HG12	1.59	0.84
8:H:91:VAL:HG12	8:H:93:SER:H	1.44	0.83
2:B:587:GLN:O	2:B:597:TYR:OH	1.97	0.81
2:B:353:LEU:HG	2:B:374:ILE:HD11	1.63	0.80
2:B:729:PRO:HD3	2:B:759:PRO:HB2	1.61	0.80
8:H:90:LEU:HG	8:H:91:VAL:HG23	1.68	0.76
10:K:51:LEU:HB3	10:K:56:LYS:HD2	1.67	0.75
4:D:184:ARG:HD2	4:D:197:GLN:HE21	1.51	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:311:LYS:HG2	2:B:323:ASN:HD21	1.51	0.73
4:D:152:LYS:HG3	4:D:157:ILE:HB	1.68	0.73
10:K:26:THR:HG23	10:K:33:SER:HA	1.70	0.73
9:I:120:SER:HA	9:I:124:LEU:HD12	1.68	0.73
2:B:357:ASP:HB2	2:B:369:ARG:HG3	1.72	0.72
9:J:248:VAL:HG12	9:J:250:GLY:H	1.55	0.71
1:A:167:G:H1	1:A:176:U:H3	1.37	0.71
1:A:142:U:H4'	4:D:268:ARG:HG3	1.71	0.70
2:B:597:TYR:H	2:B:601:ASP:HB2	1.56	0.70
9:J:87:ILE:HG22	9:J:110:LEU:HB3	1.74	0.70
3:C:68:LEU:HD12	3:C:141:LEU:HD11	1.74	0.70
3:C:148:ARG:O	3:C:152:LYS:HB2	1.92	0.69
2:B:766:SER:HB3	2:B:873:ILE:HD11	1.73	0.69
9:J:104:PHE:O	9:J:240:ASN:ND2	2.26	0.69
3:C:89:VAL:HB	3:C:119:VAL:HG22	1.74	0.68
6:F:55:ILE:HG23	6:F:94:ILE:HG12	1.77	0.67
11:L:126:GLU:OE2	11:L:129:ARG:NH2	2.25	0.67
1:A:291:G:H2'	1:A:292:G:H8	1.60	0.67
2:B:560:LEU:HD12	2:B:561:PRO:HD2	1.77	0.67
1:A:307:U:H2'	1:A:308:G:H8	1.60	0.67
3:C:40:GLN:HB2	3:C:44:TRP:CD1	2.30	0.66
1:A:31:A:O2'	1:A:264:G:OP1	2.15	0.65
2:B:295:THR:HG23	2:B:394:CYS:HB2	1.78	0.65
4:D:80:ARG:HH11	10:K:62:SER:HB3	1.61	0.65
1:A:205:U:H2'	1:A:206:G:C8	2.32	0.64
9:I:57:VAL:HG23	9:I:58:LYS:HD3	1.78	0.64
2:B:202:ILE:O	2:B:203:LYS:HG2	1.97	0.63
2:B:411:SER:HB3	2:B:575:ILE:HG13	1.80	0.63
9:J:177:VAL:HG11	9:J:214:LEU:HD13	1.79	0.63
3:C:71:ASP:HB3	3:C:74:GLU:HG2	1.79	0.63
2:B:246:ASN:O	2:B:250:SER:N	2.27	0.63
1:A:268:U:H3	1:A:305:A:H62	1.47	0.63
2:B:155:LYS:HG2	2:B:158:ARG:HH11	1.64	0.63
1:A:166:U:H2'	1:A:167:G:H8	1.64	0.62
2:B:286:LEU:HD12	2:B:581:ILE:HB	1.81	0.62
1:A:268:U:OP2	5:E:5:LYS:NZ	2.32	0.62
2:B:816:GLU:OE1	2:B:818:HIS:NE2	2.32	0.62
1:A:319:U:O4	9:J:223:LYS:NZ	2.32	0.61
8:H:124:GLU:O	8:H:128:GLN:NE2	2.32	0.61
1:A:31:A:OP1	5:E:125:LYS:NZ	2.32	0.61
4:D:22:LYS:HG3	4:D:23:PRO:HD3	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:L:51:LEU:HD11	11:L:69:LYS:HG3	1.82	0.61
2:B:297:ILE:HG12	2:B:367:ILE:HG22	1.82	0.61
2:B:371:HIS:CD2	2:B:372:PRO:HD2	2.35	0.61
1:A:205:U:H2'	1:A:206:G:H8	1.66	0.61
2:B:372:PRO:HA	2:B:375:TYR:HB3	1.82	0.61
2:B:371:HIS:HB3	2:B:374:ILE:HG22	1.83	0.61
1:A:37:A:O2'	7:G:35:VAL:O	2.18	0.60
6:F:148:ILE:HG22	6:F:150:LEU:H	1.66	0.60
4:D:77:ASP:OD1	4:D:80:ARG:NH1	2.33	0.60
2:B:802:PRO:HB2	9:I:230:SER:HB3	1.82	0.60
1:A:48:U:H2'	1:A:49:A:H8	1.67	0.60
1:A:124:U:H5'	1:A:125:U:H5'	1.84	0.60
1:A:114:C:HO2'	1:A:311:G:HO2'	1.49	0.60
2:B:370:LEU:HD13	2:B:374:ILE:HG23	1.84	0.59
2:B:581:ILE:HG13	2:B:585:GLN:HB2	1.85	0.59
3:C:126:LYS:HD2	3:C:129:MET:HE3	1.84	0.59
2:B:714:GLU:OE1	2:B:718:LYS:NZ	2.35	0.59
9:J:10:PRO:HG2	9:J:22:ALA:HB1	1.83	0.59
7:G:33:ILE:HD13	7:G:47:ARG:HD2	1.85	0.59
2:B:340:SER:HB3	2:B:353:LEU:HD13	1.85	0.59
11:L:98:PHE:HB3	11:L:101:LEU:HD21	1.85	0.59
1:A:154:C:H2'	1:A:155:A:H8	1.67	0.58
2:B:338:ILE:HG23	2:B:353:LEU:HB3	1.85	0.58
9:J:155:LYS:HZ1	9:J:232:VAL:HB	1.68	0.58
9:J:120:SER:HA	9:J:124:LEU:HB2	1.85	0.58
10:K:56:LYS:HA	10:K:59:LYS:HD2	1.85	0.58
2:B:455:HIS:NE2	2:B:533:VAL:O	2.36	0.58
1:A:193:U:H3	1:A:216:G:H1	1.51	0.57
5:E:24:GLU:OE2	9:I:28:LYS:NZ	2.37	0.57
4:D:170:LEU:HD21	9:J:249:THR:HG22	1.85	0.57
1:A:331:G:OP1	4:D:221:LYS:NZ	2.38	0.57
1:A:13:A:H2'	1:A:14:A:C8	2.39	0.57
3:C:42:LEU:HA	3:C:45:GLN:NE2	2.19	0.57
11:L:40:GLU:HG2	11:L:43:ARG:HH21	1.68	0.57
9:J:65:ILE:O	9:J:70:GLY:N	2.38	0.57
9:J:85:THR:HG22	9:J:108:ALA:HB3	1.86	0.57
4:D:266:LYS:O	4:D:268:ARG:NH1	2.38	0.57
1:A:219:G:N7	2:B:182:ARG:NH1	2.49	0.57
2:B:87:ARG:NH2	4:D:16:CYS:O	2.37	0.56
2:B:284:GLY:HA2	2:B:577:ARG:HB3	1.88	0.56
2:B:779:ALA:HB3	2:B:824:ILE:HB	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:U:H2'	1:A:331:G:H8	1.70	0.56
2:B:147:LEU:HD21	2:B:195:SER:HB2	1.86	0.56
1:A:160:G:H2'	1:A:161:G:H8	1.71	0.56
1:A:21:U:OP2	2:B:772:ARG:NH1	2.36	0.56
11:L:78:LEU:HD11	11:L:116:TYR:HB3	1.86	0.56
1:A:140:U:O2'	1:A:315:A:OP1	2.24	0.55
9:I:165:ARG:NH2	9:I:194:GLY:O	2.39	0.55
9:J:224:ALA:HA	9:J:228:LEU:HB2	1.88	0.55
1:A:147:A:N6	1:A:329:A:OP2	2.33	0.55
9:J:209:SER:HA	9:J:212:LYS:HE3	1.88	0.55
1:A:161:G:H1	1:A:182:G:H21	1.53	0.55
1:A:56:U:H3'	1:A:57:G:H8	1.72	0.55
4:D:195:GLY:O	9:J:290:ARG:NH1	2.40	0.55
2:B:801:ARG:NE	2:B:806:LYS:O	2.35	0.54
3:C:111:LEU:HD13	10:K:70:PRO:HB2	1.88	0.54
10:K:53:GLU:HA	10:K:56:LYS:HB2	1.90	0.54
2:B:287:CYS:HB3	2:B:578:MET:HE2	1.88	0.54
8:H:115:ASP:HA	8:H:118:TRP:CE2	2.43	0.54
2:B:269:PHE:O	2:B:588:GLN:NE2	2.33	0.54
9:I:137:LEU:HD13	9:I:141:LEU:HD23	1.88	0.54
11:L:14:GLU:OE2	11:L:111:ASN:ND2	2.40	0.54
4:D:139:PHE:HA	4:D:142:MET:HE2	1.89	0.54
3:C:166:PHE:HB3	3:C:169:LEU:HD13	1.90	0.54
1:A:291:G:H2'	1:A:292:G:C8	2.42	0.53
5:E:29:ALA:HB2	5:E:140:ILE:HG22	1.91	0.53
4:D:121:TYR:CZ	4:D:125:LYS:HD2	2.44	0.53
1:A:104:A:H5'	2:B:229:LYS:HB2	1.90	0.53
1:A:196:U:H2'	1:A:197:A:H8	1.73	0.53
1:A:205:U:O2'	1:A:206:G:OP1	2.26	0.53
1:A:300:A:O2'	1:A:320:U:O2'	2.26	0.53
3:C:102:LEU:HA	3:C:105:LEU:HG	1.91	0.53
4:D:28:ARG:NH1	11:L:96:GLY:O	2.39	0.53
3:C:55:LEU:HA	3:C:62:GLU:HG2	1.90	0.53
4:D:181:ALA:H	4:D:200:VAL:HG22	1.74	0.53
1:A:50:A:H2'	1:A:51:A:C8	2.44	0.53
2:B:338:ILE:N	2:B:354:GLY:O	2.42	0.52
2:B:550:ARG:NH2	2:B:555:ASP:OD2	2.36	0.52
9:I:232:VAL:HA	9:I:235:ASN:HD21	1.74	0.52
4:D:186:THR:OG1	9:J:252:GLY:O	2.24	0.52
9:I:2:LEU:HD21	9:I:204:ILE:HD12	1.90	0.52
2:B:445:THR:OG1	2:B:482:TYR:OH	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:318:THR:HG22	2:B:338:ILE:HG13	1.92	0.52
5:E:7:ARG:HA	5:E:121:GLY:HA2	1.91	0.52
2:B:296:ILE:HG13	2:B:375:TYR:CE1	2.44	0.52
2:B:288:MET:HB3	2:B:402:THR:HB	1.90	0.52
4:D:208:PHE:HB3	4:D:223:ILE:HG13	1.90	0.52
3:C:154:VAL:HA	3:C:157:ILE:HG22	1.92	0.52
4:D:233:GLU:HG2	4:D:243:ALA:HB1	1.92	0.52
9:I:107:VAL:HG23	9:I:239:ARG:HH22	1.75	0.52
1:A:14:A:H2'	1:A:15:G:C8	2.45	0.52
2:B:57:LEU:HD23	4:D:33:LEU:HD21	1.91	0.52
1:A:277:C:H4'	2:B:629:MET:HG3	1.92	0.51
1:A:305:A:OP2	5:E:3:ARG:NH1	2.43	0.51
8:H:58:LEU:HD22	8:H:119:LEU:HD11	1.91	0.51
2:B:47:PRO:HG3	2:B:66:ARG:HH12	1.75	0.51
3:C:71:ASP:O	3:C:75:ILE:HG12	2.10	0.51
5:E:108:ASP:OD2	5:E:108:ASP:N	2.43	0.51
4:D:136:TYR:HA	4:D:139:PHE:HB2	1.92	0.51
8:H:37:TRP:CD1	8:H:86:SER:HB2	2.45	0.51
11:L:11:LEU:HD21	11:L:116:TYR:HE1	1.76	0.51
3:C:50:SER:O	3:C:54:LYS:HG2	2.10	0.51
2:B:60:ASP:OD2	4:D:5:GLN:NE2	2.40	0.51
7:G:51:PHE:O	7:G:55:VAL:HG23	2.10	0.51
9:J:242:SER:OG	9:J:243:HIS:N	2.43	0.51
2:B:358:VAL:HG22	2:B:368:ILE:HG12	1.92	0.51
1:A:278:U:H2'	1:A:279:C:C6	2.46	0.51
1:A:274:U:H2'	1:A:275:U:H6	1.76	0.51
3:C:132:PHE:HD2	3:C:139:GLY:HA2	1.75	0.51
9:J:121:THR:O	9:J:152:ARG:NH2	2.44	0.51
1:A:210:G:H2'	1:A:211:G:C8	2.46	0.51
2:B:207:ASN:HD21	2:B:211:SER:HB2	1.76	0.51
10:K:30:ARG:HB2	10:K:64:CYS:HB2	1.92	0.51
1:A:19:C:H2'	1:A:20:G:H8	1.75	0.50
9:I:5:LEU:HA	9:I:42:ALA:HB3	1.92	0.50
9:J:69:PHE:HA	9:J:72:LEU:HG	1.93	0.50
1:A:25:A:N7	6:F:132:ARG:NH2	2.59	0.50
8:H:117:LEU:O	8:H:121:LYS:HG2	2.11	0.50
2:B:591:TYR:N	2:B:597:TYR:OH	2.45	0.50
10:K:23:VAL:HG23	10:K:24:LEU:H	1.76	0.50
8:H:56:SER:OG	8:H:68:ARG:NE	2.45	0.50
3:C:168:TRP:CG	10:K:22:HIS:HD1	2.30	0.50
5:E:70:SER:OG	9:J:178:ARG:NH2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:A:H2'	1:A:32:A:C8	2.47	0.50
2:B:401:VAL:HG23	2:B:558:VAL:HG13	1.94	0.50
2:B:603:PRO:HD2	2:B:604:PHE:HD1	1.77	0.50
2:B:782:TYR:OH	2:B:858:ARG:NH2	2.45	0.50
1:A:158:A:H2'	1:A:159:A:H8	1.77	0.50
9:I:3:VAL:HG22	9:I:40:HIS:HB2	1.93	0.50
10:K:41:HIS:HA	10:K:44:VAL:HG12	1.94	0.50
1:A:251:A:H2'	1:A:252:A:H8	1.76	0.49
4:D:89:LEU:HD22	10:K:35:LEU:HD11	1.94	0.49
9:J:11:GLN:HG3	9:J:16:ASP:HB2	1.94	0.49
1:A:17:A:H2'	1:A:18:U:H6	1.78	0.49
2:B:221:ALA:O	2:B:395:ARG:NH1	2.45	0.49
9:J:89:ASP:O	9:J:113:SER:OG	2.20	0.49
2:B:830:HIS:HB2	2:B:837:MET:HB2	1.94	0.49
1:A:125:U:H1'	4:D:75:LYS:HG2	1.94	0.49
1:A:164:G:H1	1:A:179:U:H3	1.61	0.49
3:C:171:ALA:O	10:K:19:ASN:ND2	2.32	0.49
6:F:31:ILE:HG21	6:F:46:TYR:HE1	1.77	0.49
4:D:214:GLY:HA2	9:J:284:LEU:HD22	1.93	0.49
1:A:66:G:H4'	1:A:67:U:O5'	2.12	0.49
1:A:307:U:H1'	2:B:94:ARG:HB2	1.94	0.49
3:C:64:TYR:HB3	3:C:65:PRO:HD3	1.94	0.49
4:D:227:GLY:N	4:D:251:ASP:OD1	2.45	0.49
9:J:121:THR:HB	9:J:149:CYS:SG	2.53	0.49
2:B:283:ASP:OD2	2:B:283:ASP:N	2.46	0.49
3:C:68:LEU:HD13	3:C:143:LEU:HB3	1.95	0.49
5:E:86:ILE:HD11	5:E:100:LEU:HD11	1.94	0.49
9:J:1:MET:HA	9:J:39:THR:HG21	1.95	0.49
2:B:158:ARG:O	2:B:162:LYS:HG2	2.13	0.49
4:D:272:ASP:OD2	4:D:272:ASP:N	2.44	0.49
9:J:126:ILE:HG13	9:J:154:VAL:HG21	1.95	0.49
2:B:653:ILE:HD12	2:B:820:LEU:HD11	1.94	0.48
3:C:172:ILE:O	10:K:15:ARG:NE	2.46	0.48
11:L:52:PHE:HD1	11:L:55:ARG:HE	1.60	0.48
11:L:101:LEU:H	11:L:101:LEU:HD23	1.78	0.48
9:I:232:VAL:HA	9:I:235:ASN:ND2	2.27	0.48
3:C:75:ILE:HG23	3:C:142:LEU:HD23	1.94	0.48
3:C:148:ARG:HH21	3:C:151:LYS:HE2	1.77	0.48
9:J:257:ASP:HA	9:J:262:ASP:HB2	1.95	0.48
1:A:275:U:C2	1:A:276:G:C8	3.01	0.48
3:C:26:PRO:HG3	3:C:173:LYS:HB3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:215:ASN:OD1	4:D:216:ILE:N	2.46	0.48
9:I:162:TYR:HA	9:I:165:ARG:HD3	1.94	0.48
3:C:88:PRO:HA	3:C:118:THR:HG23	1.94	0.48
2:B:371:HIS:HD2	2:B:372:PRO:HD2	1.77	0.48
9:I:11:GLN:HG3	9:I:16:ASP:HB3	1.95	0.48
9:J:29:THR:HA	9:J:198:PRO:HB3	1.96	0.48
10:K:71:ASN:HB3	10:K:90:LEU:HD22	1.94	0.48
2:B:237:PHE:CD2	2:B:254:LYS:HD2	2.49	0.48
2:B:348:ASP:OD1	2:B:348:ASP:N	2.46	0.48
9:J:48:ASN:HA	9:J:87:ILE:HG12	1.96	0.48
1:A:114:C:O2'	1:A:311:G:O2'	2.21	0.48
6:F:76:GLN:NE2	6:F:149:ASP:OD1	2.47	0.48
9:J:5:LEU:O	9:J:83:ARG:NH1	2.47	0.48
1:A:193:U:H2'	1:A:194:G:H8	1.78	0.47
1:A:230:U:H2'	1:A:231:A:C8	2.48	0.47
8:H:14:LEU:HD23	8:H:97:VAL:HG11	1.95	0.47
11:L:55:ARG:HD2	11:L:61:ARG:HG3	1.96	0.47
3:C:132:PHE:CD2	3:C:139:GLY:HA2	2.50	0.47
1:A:7:U:H2'	1:A:8:G:H8	1.79	0.47
4:D:36:LEU:O	11:L:24:ARG:NH1	2.47	0.47
3:C:58:ILE:HG22	3:C:61:LYS:HB2	1.97	0.47
3:C:93:VAL:HG22	3:C:140:MET:SD	2.55	0.47
4:D:231:GLN:HA	4:D:247:SER:HA	1.97	0.47
1:A:8:G:H2'	1:A:9:A:H8	1.79	0.47
1:A:188:U:O2'	1:A:189:G:OP1	2.32	0.47
1:A:282:U:H2'	1:A:283:C:C6	2.49	0.47
2:B:524:THR:OG1	2:B:525:PRO:HD3	2.14	0.47
2:B:603:PRO:HB2	2:B:728:ILE:HD11	1.96	0.47
6:F:118:PHE:CD2	6:F:137:ILE:HD11	2.50	0.47
11:L:51:LEU:HD21	11:L:69:LYS:HD3	1.96	0.47
1:A:25:A:O2'	1:A:26:A:OP2	2.23	0.47
1:A:50:A:H2'	1:A:51:A:H8	1.80	0.47
3:C:36:ARG:O	3:C:37:VAL:HG23	2.14	0.47
1:A:79:G:H22	5:E:5:LYS:HZ3	1.62	0.47
1:A:107:C:H2'	1:A:108:G:H8	1.80	0.47
1:A:109:G:N2	1:A:112:A:OP2	2.47	0.47
2:B:117:LEU:HD21	2:B:171:PRO:HD3	1.97	0.47
2:B:240:LEU:HD22	2:B:291:SER:HB2	1.97	0.47
2:B:661:TRP:CD2	2:B:728:ILE:HD13	2.50	0.47
2:B:188:LEU:O	2:B:192:ILE:HG12	2.15	0.46
9:I:149:CYS:HA	9:I:152:ARG:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:L:74:LEU:HD21	11:L:119:LEU:HD11	1.97	0.46
2:B:234:GLN:HB3	2:B:237:PHE:O	2.15	0.46
6:F:49:ILE:HD13	6:F:94:ILE:HD12	1.96	0.46
1:A:92:C:H42	1:A:308:G:H1	1.63	0.46
1:A:302:C:OP1	9:J:187:ARG:NH2	2.49	0.46
2:B:560:LEU:HD23	2:B:565:LEU:HA	1.96	0.46
4:D:115:LYS:HB2	4:D:215:ASN:O	2.16	0.46
4:D:277:ILE:HD11	9:J:234:LEU:HD12	1.97	0.46
7:G:60:SER:OG	7:G:61:SER:N	2.49	0.46
2:B:566:LEU:HB3	2:B:567:PRO:HD3	1.98	0.46
2:B:662:ARG:HD3	2:B:665:GLN:HE21	1.80	0.46
2:B:802:PRO:O	6:F:126:ASN:ND2	2.49	0.46
1:A:318:A:O2'	1:A:319:U:OP1	2.32	0.46
2:B:664:LEU:HB2	2:B:823:PHE:HE2	1.81	0.46
1:A:157:U:OP2	2:B:111:ARG:NH1	2.46	0.46
2:B:92:LEU:HB2	11:L:95:LEU:HD13	1.98	0.46
4:D:11:CYS:O	4:D:14:THR:HG22	2.15	0.46
6:F:149:ASP:OD1	6:F:149:ASP:N	2.48	0.46
2:B:859:ASN:HB2	2:B:862:THR:HB	1.98	0.46
3:C:91:LEU:HA	3:C:142:LEU:HA	1.98	0.46
3:C:95:ASN:HB3	3:C:96:LYS:H	1.58	0.46
6:F:57:LYS:O	6:F:61:LYS:HG3	2.15	0.46
9:J:139:THR:OG1	9:J:140:PHE:N	2.49	0.46
9:J:7:VAL:HB	9:J:43:ILE:HA	1.97	0.45
3:C:91:LEU:HD12	3:C:142:LEU:HB3	1.99	0.45
9:J:87:ILE:HG13	9:J:87:ILE:O	2.16	0.45
2:B:190:ARG:HA	2:B:193:LYS:HG2	1.98	0.45
1:A:7:U:H2'	1:A:8:G:C8	2.52	0.45
1:A:75:U:H5''	2:B:772:ARG:HA	1.98	0.45
1:A:207:C:H2'	1:A:208:A:H8	1.81	0.45
2:B:732:LEU:HD23	2:B:732:LEU:H	1.81	0.45
6:F:35:ALA:HB1	6:F:150:LEU:HD22	1.99	0.45
9:I:86:LEU:O	9:I:110:LEU:N	2.44	0.45
1:A:25:A:H4'	1:A:26:A:O5'	2.17	0.45
1:A:331:G:H2'	1:A:332:G:H8	1.81	0.45
2:B:442:GLN:HA	2:B:549:ARG:HH11	1.81	0.45
9:I:19:THR:HG22	9:I:21:GLN:H	1.81	0.45
4:D:175:MET:O	10:K:9:TYR:OH	2.35	0.45
9:I:49:HIS:CD2	9:I:88:ILE:HG12	2.52	0.45
9:J:48:ASN:ND2	9:J:50:SER:H	2.15	0.45
3:C:26:PRO:HB3	3:C:29:ASN:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:6:ASN:ND2	9:I:159:VAL:HG22	2.31	0.44
10:K:18:TYR:O	10:K:22:HIS:HD2	1.99	0.44
1:A:190:A:H1'	2:B:161:SER:OG	2.17	0.44
1:A:227:C:O2'	1:A:228:G:OP1	2.33	0.44
2:B:450:ALA:HB3	2:B:545:LEU:HD23	2.00	0.44
6:F:151:ASN:OD1	6:F:153:HIS:ND1	2.50	0.44
9:I:30:LEU:HD22	9:I:41:ILE:HG21	2.00	0.44
9:I:72:LEU:O	9:I:76:THR:HG22	2.17	0.44
11:L:109:LEU:HA	11:L:112:VAL:HG12	1.99	0.44
1:A:11:C:H2'	1:A:12:A:H8	1.82	0.44
1:A:11:C:C2	1:A:12:A:C8	3.05	0.44
3:C:113:TYR:HD2	3:C:164:LEU:HG	1.81	0.44
9:I:65:ILE:HG21	9:I:80:LEU:HB2	1.99	0.44
1:A:45:G:N7	6:F:90:LYS:NZ	2.52	0.44
2:B:447:ALA:HB3	2:B:487:ILE:HG23	1.98	0.44
3:C:40:GLN:HB2	3:C:44:TRP:HD1	1.78	0.44
2:B:740:PHE:CZ	2:B:845:HIS:HB2	2.52	0.44
6:F:27:LYS:HE3	6:F:46:TYR:HB2	2.00	0.44
11:L:55:ARG:NH2	11:L:56:ARG:HB2	2.32	0.44
1:A:265:U:H4'	1:A:266:U:H5'	1.99	0.44
2:B:455:HIS:CE1	2:B:533:VAL:HG23	2.53	0.44
7:G:25:THR:OG1	7:G:26:HIS:N	2.50	0.44
9:I:159:VAL:HG11	9:I:162:TYR:HD2	1.83	0.44
9:I:212:LYS:HA	9:I:216:LEU:O	2.17	0.44
1:A:19:C:H2'	1:A:20:G:C8	2.53	0.44
3:C:58:ILE:CG2	3:C:61:LYS:HE2	2.48	0.44
4:D:187:LYS:HB2	4:D:229:VAL:HB	1.99	0.44
4:D:238:ASP:OD1	4:D:238:ASP:N	2.50	0.44
9:J:48:ASN:HB3	9:J:51:GLU:OE2	2.18	0.44
1:A:15:G:H2'	1:A:83:G:N7	2.32	0.44
4:D:118:LEU:O	4:D:122:ILE:HG12	2.17	0.44
4:D:172:LYS:O	4:D:176:ALA:N	2.50	0.44
9:I:65:ILE:HG23	9:I:69:PHE:HB2	2.00	0.44
2:B:309:PHE:HD2	2:B:389:LEU:HD13	1.83	0.43
4:D:208:PHE:HD1	4:D:225:LYS:HE2	1.83	0.43
3:C:41:PRO:HA	3:C:44:TRP:HB2	1.99	0.43
3:C:94:CYS:SG	3:C:139:GLY:HA3	2.58	0.43
3:C:166:PHE:CD2	3:C:168:TRP:HB2	2.54	0.43
4:D:106:LEU:O	4:D:110:ASN:HB2	2.18	0.43
5:E:56:LEU:HD12	5:E:60:LEU:HD23	2.00	0.43
7:G:24:LEU:HG	7:G:29:ILE:HB	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:A:H2'	1:A:17:A:C8	2.53	0.43
3:C:146:ASP:HA	3:C:151:LYS:NZ	2.34	0.43
1:A:37:A:O4'	7:G:34:PHE:HB3	2.19	0.43
1:A:218:U:H4'	1:A:219:G:H5'	1.99	0.43
2:B:205:LEU:HD23	2:B:205:LEU:H	1.83	0.43
2:B:502:GLN:O	2:B:505:ASN:HB2	2.18	0.43
9:I:11:GLN:NE2	9:I:16:ASP:O	2.51	0.43
1:A:230:U:OP2	11:L:43:ARG:HD2	2.18	0.43
9:I:160:TYR:CZ	9:I:193:SER:HB2	2.54	0.43
1:A:254:C:H2'	1:A:255:U:H6	1.84	0.43
4:D:154:LEU:HD11	10:K:9:TYR:HD1	1.83	0.43
1:A:96:C:O2'	1:A:97:A:OP1	2.32	0.43
1:A:181:U:H3'	1:A:182:G:C8	2.53	0.43
3:C:175:ARG:HH21	3:C:177:THR:HG1	1.66	0.43
4:D:152:LYS:HA	4:D:157:ILE:HD12	2.00	0.43
4:D:16:CYS:HB3	11:L:94:ALA:HB2	2.01	0.43
4:D:70:ASP:OD1	4:D:71:LYS:N	2.52	0.43
1:A:75:U:H2'	1:A:76:U:C6	2.54	0.43
1:A:159:A:O3'	2:B:235:LYS:NZ	2.50	0.43
1:A:274:U:H2'	1:A:275:U:C6	2.53	0.43
2:B:234:GLN:C	2:B:236:HIS:H	2.26	0.43
3:C:52:ILE:O	3:C:56:ILE:HG12	2.19	0.43
1:A:152:C:H2'	1:A:153:U:C6	2.54	0.43
2:B:166:MET:O	2:B:170:MET:HG2	2.19	0.43
2:B:362:GLN:HG3	2:B:488:ASN:HD21	1.84	0.43
2:B:526:ASN:HA	11:L:72:ARG:HH12	1.83	0.43
2:B:714:GLU:O	2:B:718:LYS:HG3	2.19	0.43
4:D:81:ILE:HD11	10:K:62:SER:HA	2.00	0.43
9:I:87:ILE:HA	9:I:110:LEU:HB3	2.01	0.42
10:K:55:LEU:O	10:K:59:LYS:HB3	2.19	0.42
2:B:334:PHE:CZ	2:B:358:VAL:HB	2.54	0.42
3:C:93:VAL:HG21	3:C:108:ILE:HD12	2.01	0.42
4:D:200:VAL:HG12	4:D:210:MET:HG2	2.01	0.42
9:J:89:ASP:OD1	9:J:89:ASP:N	2.48	0.42
10:K:60:PHE:CE1	10:K:67:VAL:HG22	2.53	0.42
10:K:103:GLN:HA	10:K:114:LEU:HA	1.99	0.42
1:A:114:C:H2'	1:A:115:G:C8	2.54	0.42
1:A:198:C:H2'	1:A:199:C:H6	1.85	0.42
2:B:203:LYS:HE2	2:B:203:LYS:HA	2.01	0.42
2:B:362:GLN:HB3	2:B:364:ASP:OD1	2.19	0.42
2:B:700:ILE:HG23	2:B:705:ASP:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:47:LEU:O	3:C:51:ILE:HG13	2.18	0.42
1:A:201:C:H42	1:A:208:A:H61	1.67	0.42
2:B:712:ASP:OD2	2:B:713:TYR:N	2.53	0.42
2:B:272:THR:HB	2:B:589:ILE:HD11	2.00	0.42
2:B:678:ASP:OD1	2:B:679:LYS:N	2.53	0.42
5:E:105:LYS:HD2	5:E:110:ASP:OD1	2.19	0.42
9:J:34:HIS:NE2	9:J:76:THR:O	2.52	0.42
11:L:113:ARG:NH1	11:L:117:MET:HE1	2.34	0.42
1:A:26:A:H2'	1:A:27:A:C8	2.54	0.42
1:A:161:G:H1	1:A:182:G:N2	2.17	0.42
1:A:262:U:P	2:B:269:PHE:HB3	2.60	0.42
2:B:273:HIS:CD2	2:B:831:LEU:HD23	2.55	0.42
6:F:64:LYS:O	6:F:68:MET:HG2	2.19	0.42
10:K:101:CYS:SG	10:K:116:GLU:HA	2.60	0.42
4:D:99:ASN:ND2	4:D:132:SER:OG	2.53	0.42
3:C:51:ILE:HG22	3:C:55:LEU:HD23	2.01	0.41
8:H:14:LEU:HD22	8:H:67:ILE:HD12	2.02	0.41
1:A:206:G:H2'	1:A:207:C:O4'	2.20	0.41
3:C:61:LYS:HA	3:C:64:TYR:HB2	2.02	0.41
4:D:197:GLN:O	4:D:212:VAL:HG12	2.20	0.41
4:D:4:THR:O	4:D:8:ILE:HG12	2.20	0.41
1:A:160:G:H2'	1:A:161:G:C8	2.53	0.41
4:D:154:LEU:HD12	10:K:12:ARG:HD3	2.02	0.41
9:J:9:TRP:HE1	9:J:18:VAL:HG21	1.85	0.41
1:A:8:G:H2'	1:A:9:A:C8	2.55	0.41
2:B:429:LYS:HG3	2:B:462:LEU:HD23	2.02	0.41
3:C:44:TRP:CZ3	3:C:47:LEU:HD13	2.55	0.41
9:J:65:ILE:HA	9:J:69:PHE:HD2	1.85	0.41
3:C:154:VAL:O	3:C:158:GLN:HG2	2.21	0.41
4:D:269:ARG:HG2	4:D:271:ASP:HB2	2.02	0.41
9:I:224:ALA:O	9:I:228:LEU:HB2	2.20	0.41
2:B:274:ARG:HA	2:B:279:THR:HG21	2.02	0.41
4:D:230:PHE:HB2	4:D:248:ILE:HG12	2.02	0.41
8:H:61:ASP:HB2	8:H:64:LEU:HB3	2.03	0.41
1:A:101:C:O2'	1:A:102:G:OP1	2.34	0.41
11:L:93:ILE:HG13	11:L:99:VAL:HA	2.02	0.41
1:A:158:A:H2'	1:A:159:A:C8	2.54	0.41
1:A:163:G:H1	1:A:180:C:H42	1.69	0.41
1:A:229:G:H2'	1:A:230:U:C6	2.56	0.41
2:B:332:VAL:HG21	2:B:442:GLN:HG2	2.02	0.41
2:B:359:PHE:HB3	2:B:367:ILE:HG12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:642:HIS:NE2	2:B:653:ILE:HD13	2.36	0.41
4:D:2:ASP:HB3	4:D:5:GLN:HE22	1.85	0.41
5:E:16:PRO:HG2	9:I:199:LEU:HD23	2.03	0.41
5:E:101:MET:HE1	9:I:210:LEU:HB2	2.03	0.41
9:I:7:VAL:HB	9:I:43:ILE:HA	2.02	0.41
2:B:85:SER:O	2:B:85:SER:OG	2.38	0.40
2:B:295:THR:HA	2:B:368:ILE:O	2.21	0.40
3:C:108:ILE:HG13	3:C:109:PRO:HD3	2.02	0.40
3:C:131:THR:O	3:C:135:VAL:HG22	2.21	0.40
4:D:143:TYR:OH	10:K:25:PRO:HG2	2.21	0.40
9:J:207:VAL:O	9:J:211:ILE:HG12	2.21	0.40
1:A:77:U:H2'	1:A:78:G:H8	1.86	0.40
1:A:167:G:O6	1:A:177:A:N6	2.55	0.40
1:A:180:C:H2'	1:A:181:U:H6	1.86	0.40
2:B:493:GLU:O	2:B:499:SER:OG	2.40	0.40
1:A:309:G:H2'	1:A:310:A:H8	1.86	0.40
3:C:34:TRP:HB2	3:C:35:PRO:HD3	2.02	0.40
5:E:26:VAL:O	9:I:75:ARG:HD3	2.21	0.40
9:I:133:TYR:HB3	9:I:173:PHE:HD1	1.86	0.40
9:J:242:SER:OG	9:J:247:ILE:O	2.30	0.40
10:K:60:PHE:HE1	10:K:67:VAL:HG22	1.86	0.40
1:A:104:A:N7	2:B:155:LYS:NZ	2.66	0.40
1:A:145:A:H4'	1:A:150:A:N1	2.36	0.40
1:A:212:U:H2'	1:A:213:A:H8	1.86	0.40
2:B:234:GLN:HG2	2:B:239:TRP:CE2	2.57	0.40
5:E:12:GLU:OE1	5:E:133:ARG:NH2	2.54	0.40
5:E:16:PRO:HA	5:E:17:PRO:HD3	1.89	0.40
9:I:158:ILE:HD13	9:I:177:VAL:HG23	2.04	0.40
9:J:160:TYR:CZ	9:J:207:VAL:HG22	2.56	0.40
1:A:183:C:H2'	1:A:184:U:C6	2.57	0.40
1:A:218:U:OP2	2:B:182:ARG:NH2	2.54	0.40
2:B:536:LYS:O	2:B:540:ASP:HB2	2.22	0.40
2:B:553:THR:HG22	2:B:590:GLN:NE2	2.36	0.40
3:C:58:ILE:HD12	3:C:58:ILE:HA	1.95	0.40
6:F:83:SER:HB3	6:F:91:MET:SD	2.61	0.40
8:H:13:LYS:HG2	8:H:100:LEU:HD12	2.04	0.40
9:J:129:LEU:HD12	9:J:129:LEU:HA	1.89	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	
2	B	785/875 (90%)	730 (93%)	54 (7%)	1 (0%)	48	68
3	C	150/195 (77%)	129 (86%)	21 (14%)	0	100	100
4	D	247/279 (88%)	240 (97%)	7 (3%)	0	100	100
5	E	143/173 (83%)	138 (96%)	5 (4%)	0	100	100
6	F	155/158 (98%)	152 (98%)	3 (2%)	0	100	100
7	G	117/140 (84%)	114 (97%)	3 (3%)	0	100	100
8	H	129/133 (97%)	124 (96%)	5 (4%)	0	100	100
9	I	240/293 (82%)	221 (92%)	18 (8%)	1 (0%)	30	49
9	J	290/293 (99%)	277 (96%)	13 (4%)	0	100	100
10	K	100/198 (50%)	89 (89%)	10 (10%)	1 (1%)	12	24
11	L	129/201 (64%)	126 (98%)	3 (2%)	0	100	100
All	All	2485/2938 (85%)	2340 (94%)	142 (6%)	3 (0%)	49	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	593	ASN
10	K	23	VAL
9	I	138	PRO

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	
2	B	718/785 (92%)	718 (100%)	0	100	100
3	C	144/185 (78%)	143 (99%)	1 (1%)	76	89
4	D	235/261 (90%)	235 (100%)	0	100	100
5	E	133/160 (83%)	133 (100%)	0	100	100
6	F	148/149 (99%)	148 (100%)	0	100	100
7	G	110/127 (87%)	110 (100%)	0	100	100
8	H	122/123 (99%)	122 (100%)	0	100	100
9	I	216/258 (84%)	216 (100%)	0	100	100
9	J	257/258 (100%)	257 (100%)	0	100	100
10	K	95/184 (52%)	95 (100%)	0	100	100
11	L	118/182 (65%)	117 (99%)	1 (1%)	73	88
All	All	2296/2672 (86%)	2294 (100%)	2 (0%)	87	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	37	VAL
11	L	62	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	61	GLN
2	B	78	HIS
2	B	180	HIS
2	B	201	ASN
2	B	246	ASN
2	B	273	HIS
2	B	323	ASN
2	B	342	ASN
2	B	380	ASN
2	B	488	ASN
2	B	570	HIS
2	B	573	ASN
2	B	585	GLN
2	B	590	GLN
2	B	669	ASN
2	B	677	ASN

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Mol	Chain	Res	Type
2	B	774	HIS
2	B	850	GLN
3	C	45	GLN
3	C	77	GLN
3	C	107	GLN
3	C	156	GLN
4	D	99	ASN
4	D	135	GLN
4	D	156	ASN
4	D	179	ASN
4	D	197	GLN
6	F	77	GLN
6	F	111	GLN
6	F	113	ASN
7	G	28	GLN
7	G	30	HIS
8	H	39	ASN
8	H	71	HIS
8	H	128	GLN
9	I	34	HIS
9	I	49	HIS
9	I	123	ASN
9	I	168	GLN
9	J	11	GLN
9	J	48	ASN
9	J	49	HIS
9	J	245	GLN
11	L	47	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	329/340 (96%)	79 (24%)	13 (3%)

All (79) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	16	A
1	A	25	A
1	A	26	A
1	A	30	G

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Mol	Chain	Res	Type
1	A	36	U
1	A	40	A
1	A	41	A
1	A	44	G
1	A	45	G
1	A	54	U
1	A	55	U
1	A	56	U
1	A	66	G
1	A	67	U
1	A	68	A
1	A	69	A
1	A	71	A
1	A	72	U
1	A	79	G
1	A	81	U
1	A	82	U
1	A	83	G
1	A	84	A
1	A	86	A
1	A	88	U
1	A	89	C
1	A	94	A
1	A	97	A
1	A	98	A
1	A	102	G
1	A	105	U
1	A	113	A
1	A	114	C
1	A	123	A
1	A	124	U
1	A	125	U
1	A	126	U
1	A	143	U
1	A	144	U
1	A	147	A
1	A	171	U
1	A	172	U
1	A	186	C
1	A	187	U
1	A	189	G
1	A	190	A

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Mol	Chain	Res	Type
1	A	203	A
1	A	206	G
1	A	218	U
1	A	219	G
1	A	220	U
1	A	224	C
1	A	225	U
1	A	228	G
1	A	233	U
1	A	234	G
1	A	243	U
1	A	244	U
1	A	248	U
1	A	262	U
1	A	263	A
1	A	266	U
1	A	267	A
1	A	269	G
1	A	271	C
1	A	281	U
1	A	286	G
1	A	290	U
1	A	295	C
1	A	303	C
1	A	305	A
1	A	306	A
1	A	316	C
1	A	318	A
1	A	319	U
1	A	321	C
1	A	322	U
1	A	328	C
1	A	329	A

All (13) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	15	G
1	A	24	C
1	A	25	A
1	A	66	G
1	A	82	U

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Mol	Chain	Res	Type
1	A	96	C
1	A	101	C
1	A	104	A
1	A	143	U
1	A	188	U
1	A	205	U
1	A	227	C
1	A	318	A

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, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

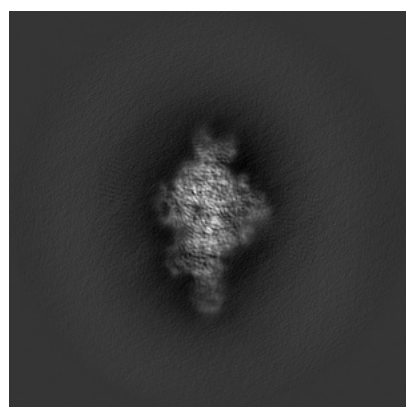
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30296. These allow visual inspection of the internal detail of the map and identification of artifacts.

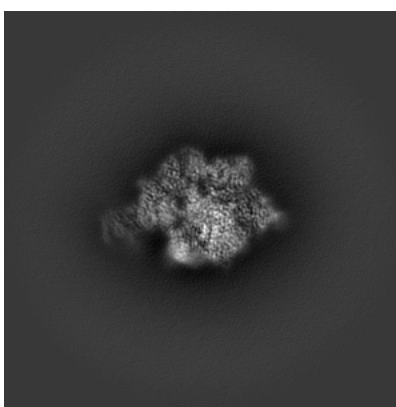
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

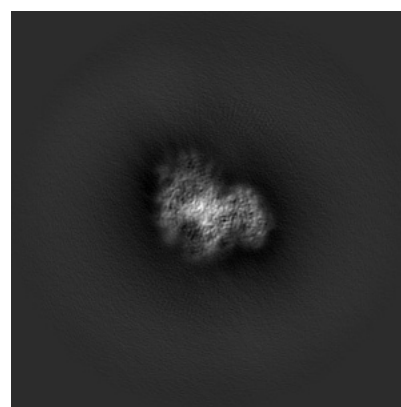
6.1.1 Primary map



X



Y

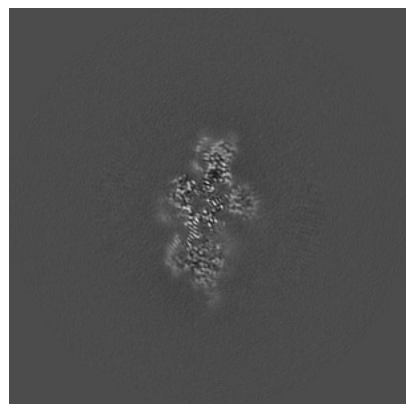


Z

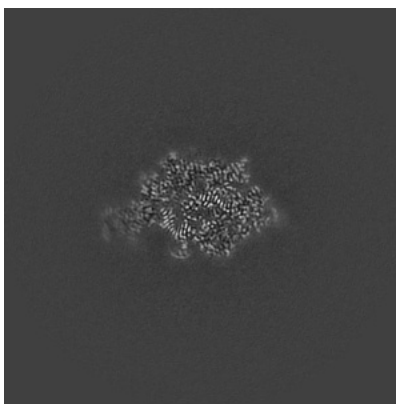
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

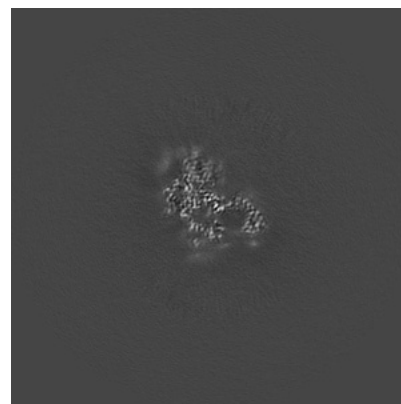
6.2.1 Primary map



X Index: 192



Y Index: 192

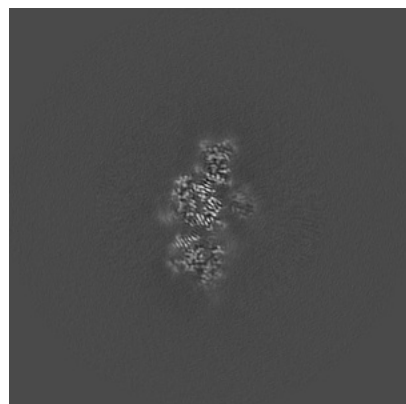


Z Index: 192

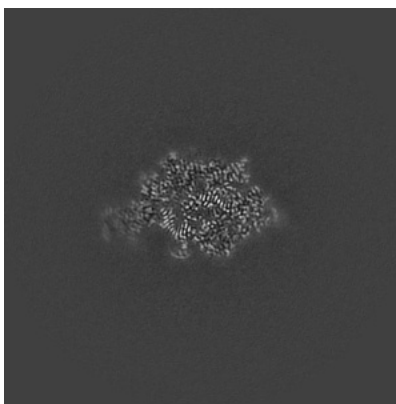
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

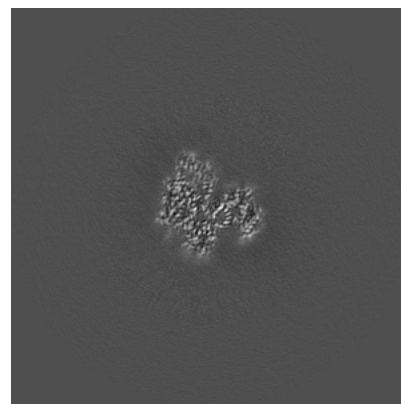
6.3.1 Primary map



X Index: 195



Y Index: 192

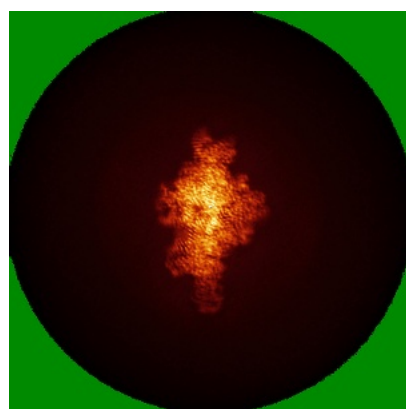


Z Index: 203

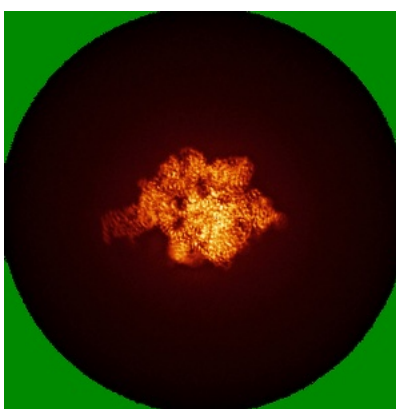
The images above show the largest variance slices of the map in three orthogonal directions.

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

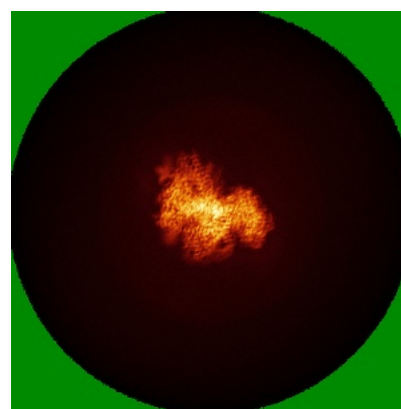
6.4.1 Primary map



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

This section was not generated.

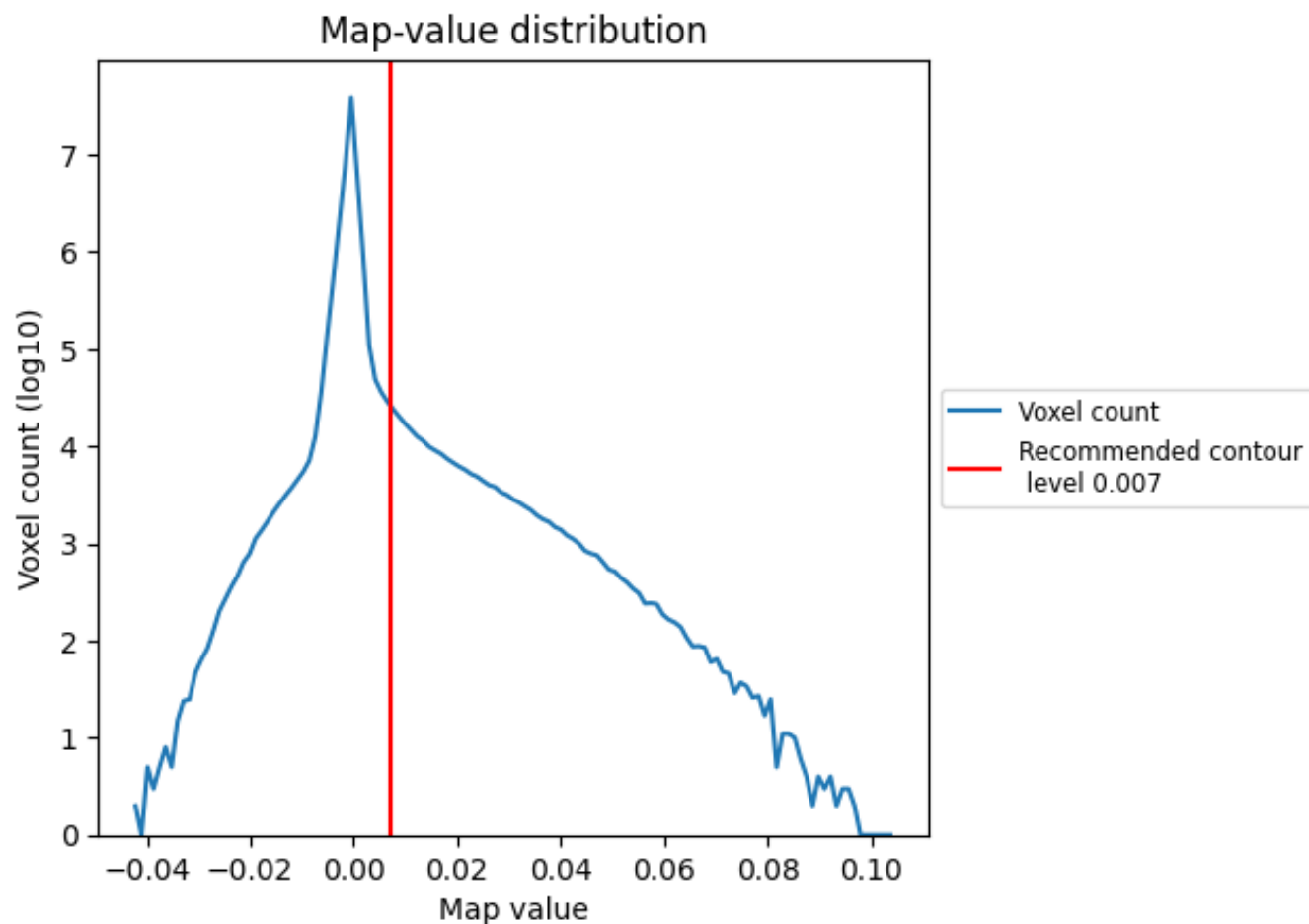
6.6 Mask visualisation

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

7 Map analysis [i](#)

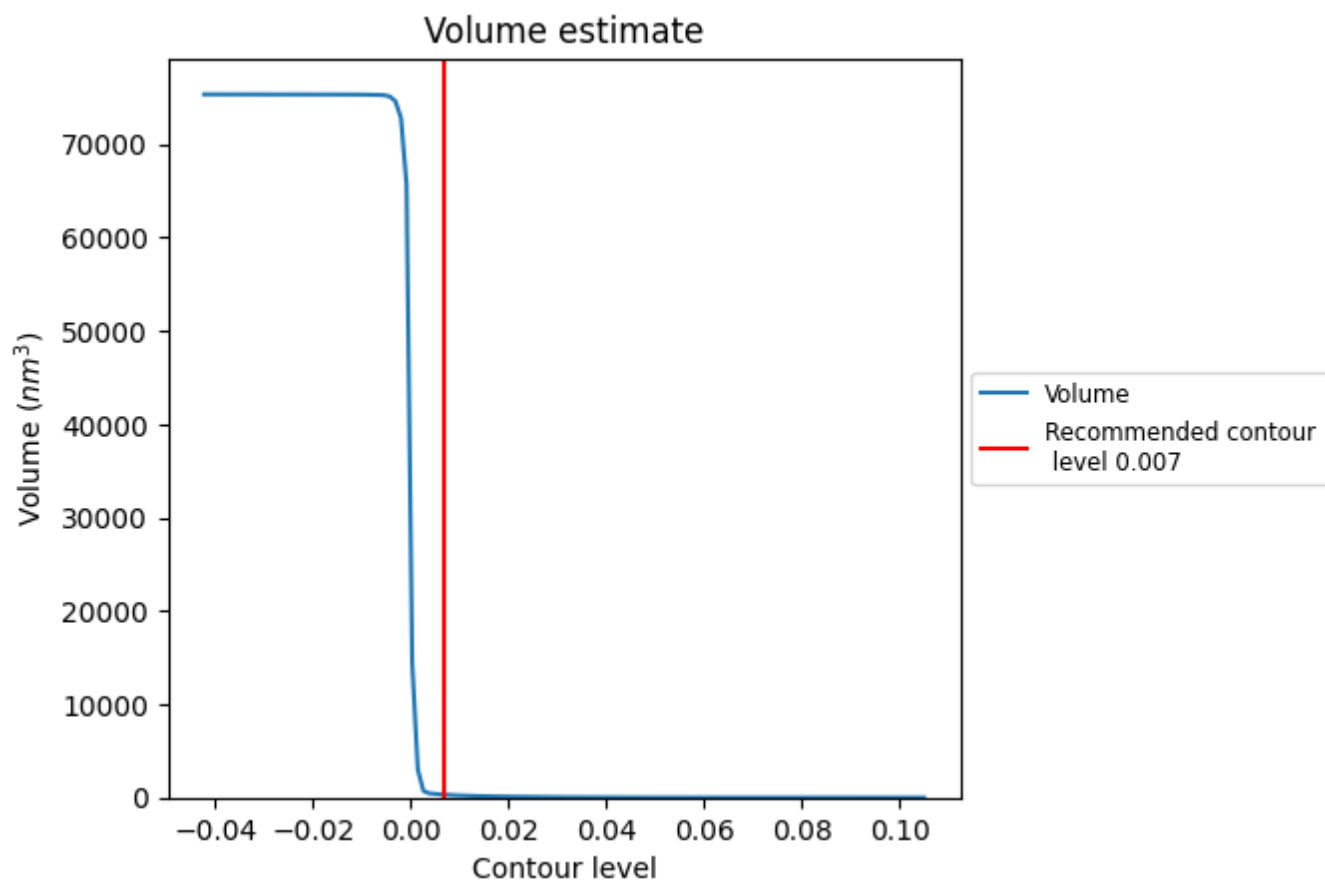
This section contains the results of statistical analysis of the map.

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.

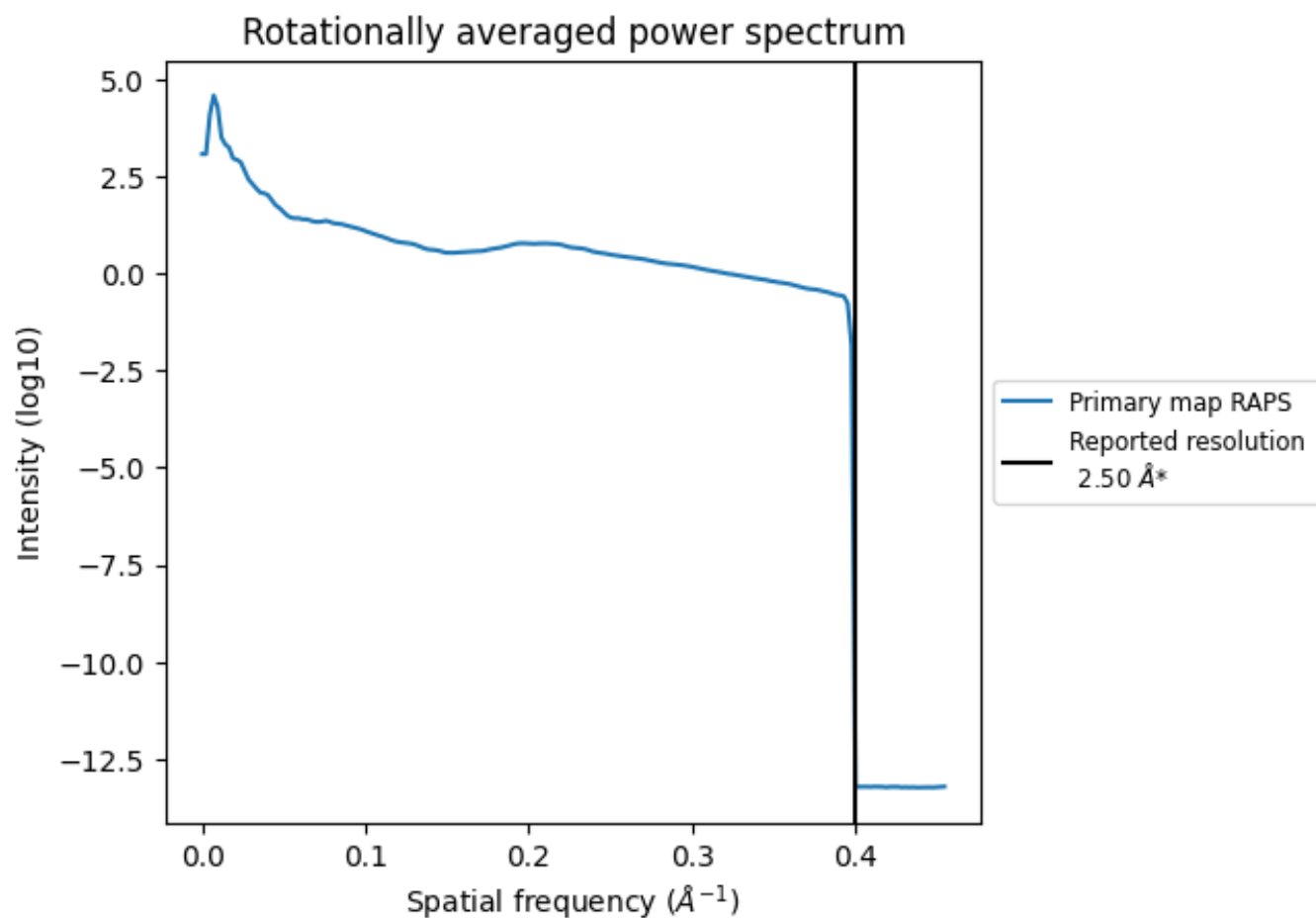
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 302 nm³; this corresponds to an approximate mass of 273 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 0.400 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

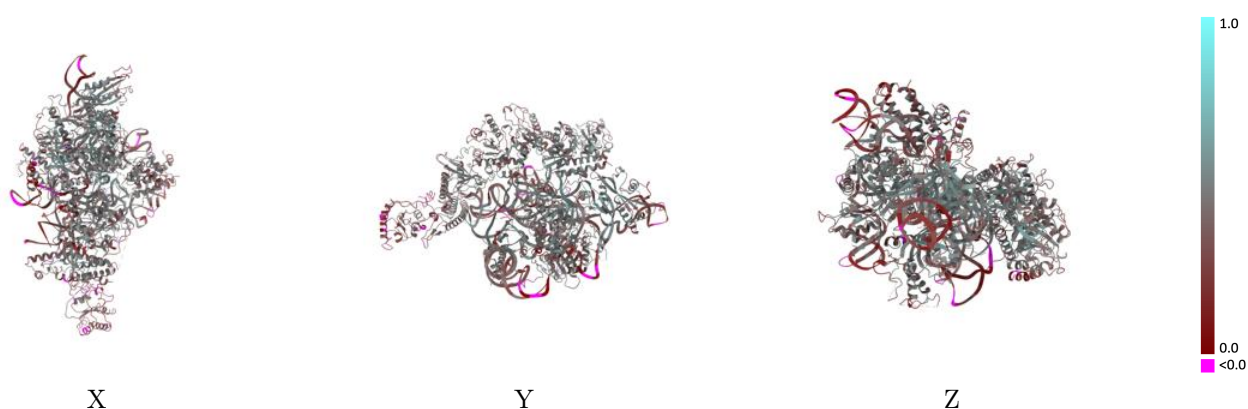
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30296 and PDB model 7C79. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

This section was not generated.

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

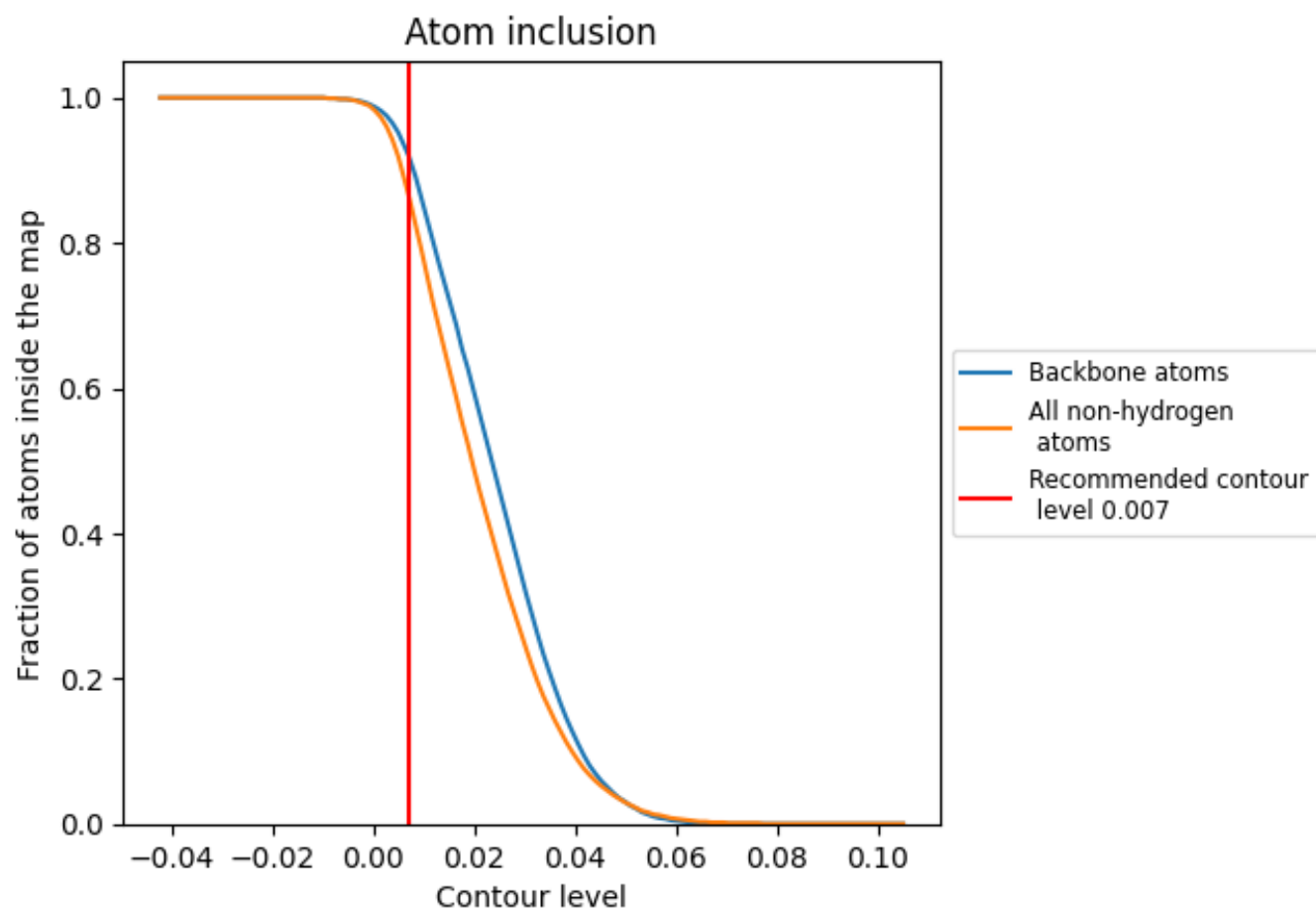


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion ⓘ



At the recommended contour level, 92% of all backbone atoms, 86% 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 (0.007) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8650	0.4020
A	0.8770	0.3610
B	0.8920	0.4380
C	0.6030	0.2430
D	0.8220	0.4020
E	0.9120	0.4620
F	0.9020	0.4350
G	0.9200	0.4830
H	0.8080	0.3590
I	0.8980	0.4570
J	0.8900	0.4140
K	0.8480	0.3720
L	0.8390	0.3970

1.0

0.0

<0.0