



Full wwPDB EM Validation Report ⓘ

Nov 9, 2022 – 02:26 AM JST

PDB ID : 6AGB
EMDB ID : EMD-9616
Title : Cryo-EM structure of yeast Ribonuclease P
Authors : Lan, P.; Tan, M.; Wu, J.; Lei, M.
Deposited on : 2018-08-10
Resolution : 3.48 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) : 2.31.2

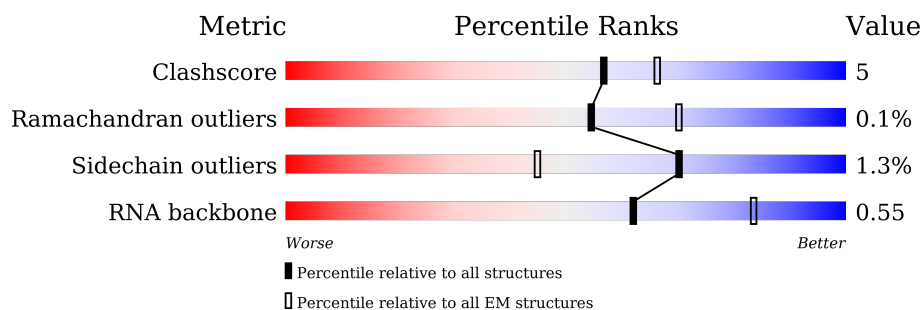
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.48 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
RNA backbone	4643	859

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	369	
2	B	875	
3	C	195	
4	D	279	
5	E	173	
6	F	158	
7	G	140	

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Mol	Chain	Length	Quality of chain
8	H	133	<div><div></div><div>14%</div><div>87%</div><div>11%</div><div></div></div>
9	I	293	<div><div></div><div></div><div>70%</div><div>13%</div><div>17%</div><div></div></div>
9	J	293	<div><div></div><div></div><div>84%</div><div>16%</div><div></div></div>
10	K	144	<div><div></div><div>9%</div><div>70%</div><div>18%</div><div>11%</div><div></div></div>

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 27190 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 P RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	369	Total	C	N	O	P	0	0
			7861	3509	1387	2596	369		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	784	Total	C	N	O	S	0	0
			6389	4046	1161	1146	36		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	175	Total	C	N	O	S	0	0
			1435	935	241	250	9		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	227	Total	C	N	O	S	0	0
			1886	1216	316	343	11		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	146	Total	C	N	O	S	0	0
			1141	719	202	212	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
			1272	804	222	242	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
			961	609	167	183	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	293	Total	C	N	O	S	0	0
			2260	1415	413	422	10		

- Molecule 10 is a protein called Ribonuclease P protein subunit RPR2.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	K	128	Total	C	N	O	S	0	0
			1024	642	193	183	6		

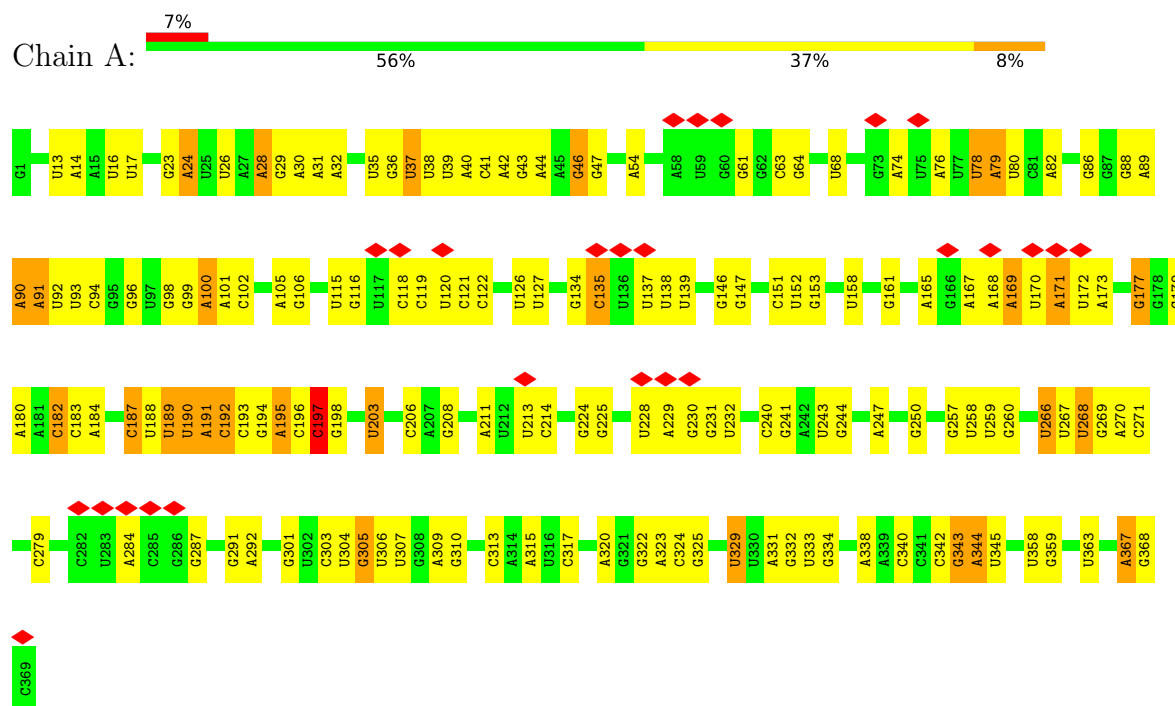
- Molecule 11 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

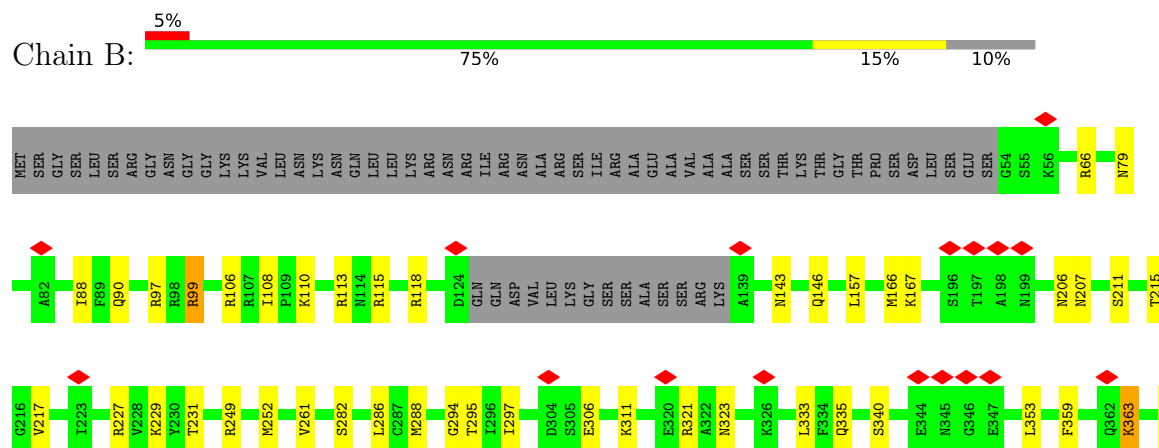
3 Residue-property plots

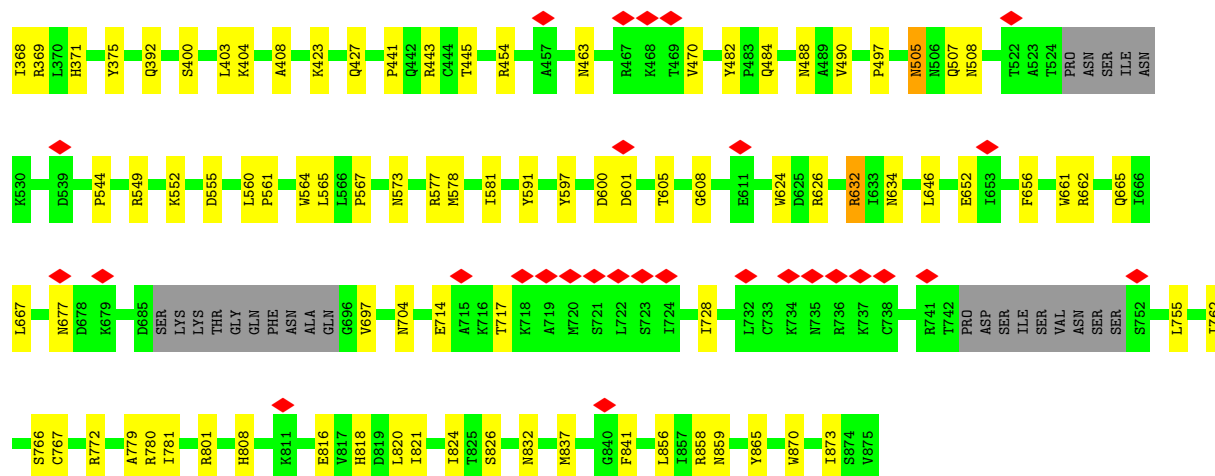
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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 P RNA

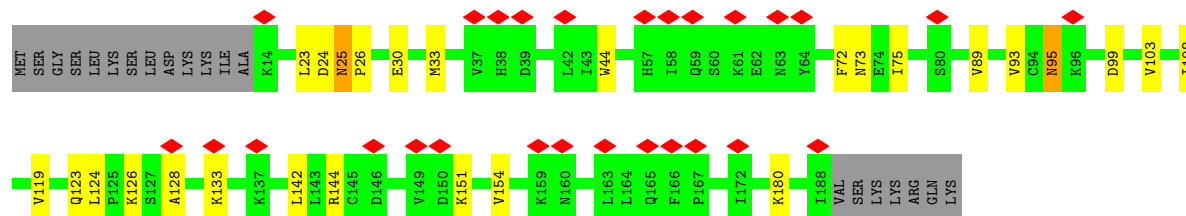
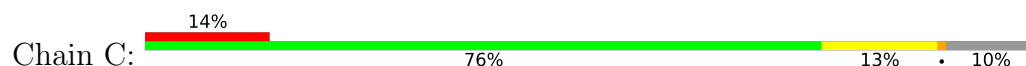


• Molecule 2: Ribonucleases P/MRP protein subunit POP1

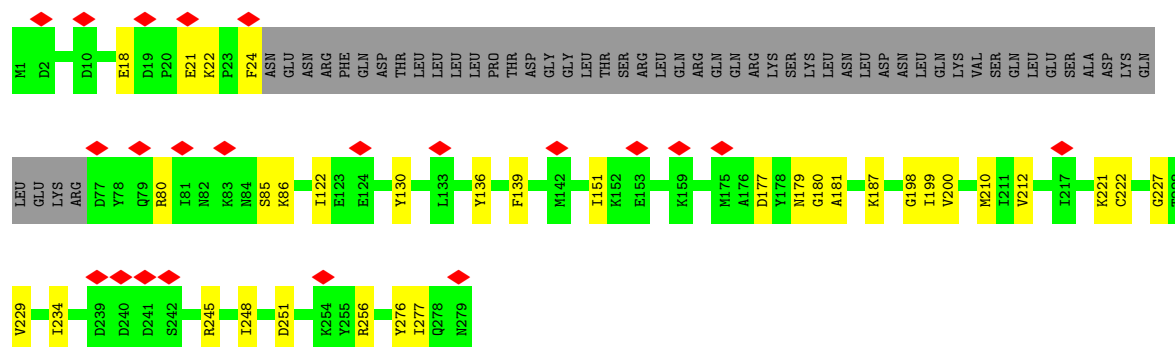




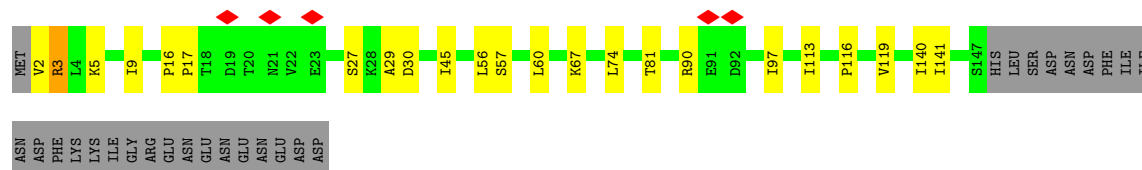
• Molecule 3: Ribonucleases P/MRP protein subunit POP3



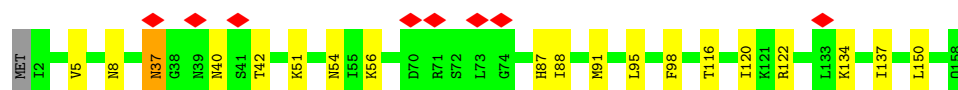
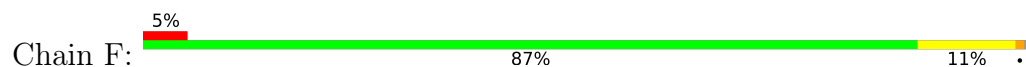
• Molecule 4: RNases MRP/P 32.9 kDa subunit



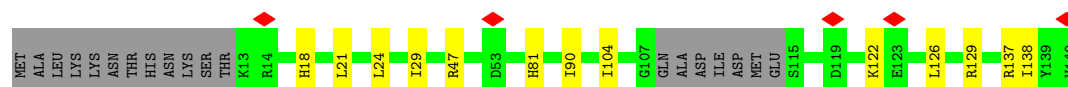
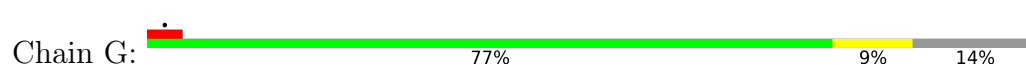
• Molecule 5: Ribonuclease P/MRP protein subunit POP5



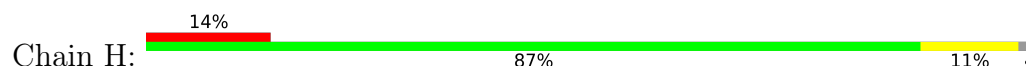
- Molecule 6: Ribonucleases P/MRP protein subunit POP6



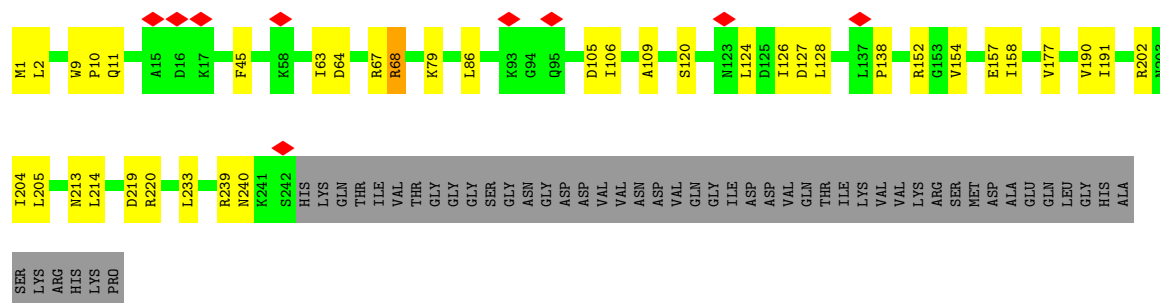
- Molecule 7: Ribonucleases P/MRP protein subunit POP7



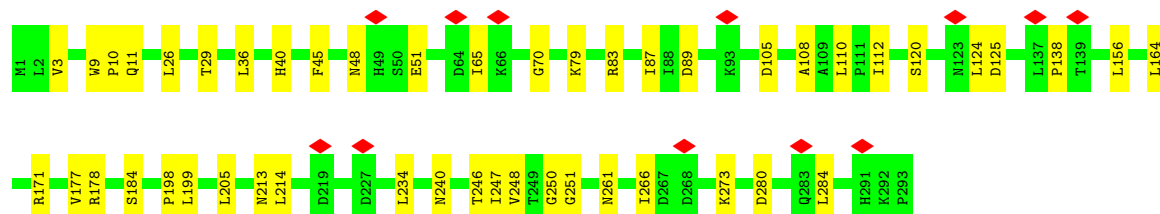
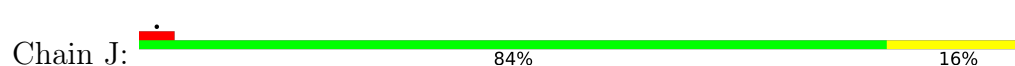
- Molecule 8: Ribonucleases P/MRP protein subunit POP8



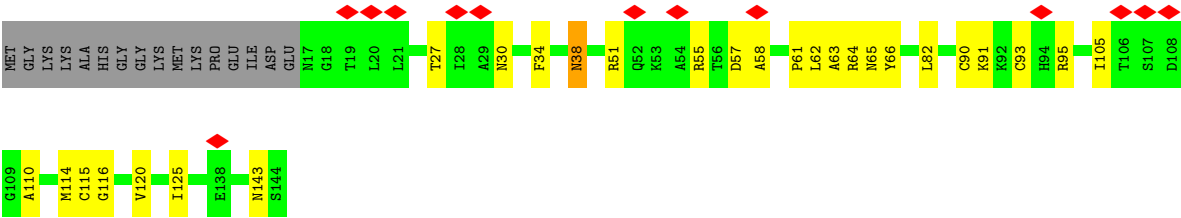
- Molecule 9: Ribonuclease P/MRP protein subunit RPP1



- Molecule 9: Ribonuclease P/MRP protein subunit RPP1



- Molecule 10: Ribonuclease P protein subunit RPR2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	164765	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 ($e^-/\text{\AA}^2$)	50.0	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.291	Depositor
Minimum map value	-0.138	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	337.92, 337.92, 337.92	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.32, 1.32, 1.32	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

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	0.23	0/8786	0.89	15/13692 (0.1%)
2	B	0.26	0/6527	0.50	0/8792
3	C	0.28	0/1473	0.52	0/2005
4	D	0.25	0/1921	0.47	0/2578
5	E	0.25	0/1154	0.47	0/1550
6	F	0.26	0/1290	0.48	0/1733
7	G	0.24	0/973	0.48	0/1305
8	H	0.26	0/1101	0.52	1/1488 (0.1%)
9	I	0.25	0/1908	0.50	0/2577
9	J	0.24	0/2291	0.52	1/3092 (0.0%)
10	K	0.29	0/1044	0.53	0/1410
All	All	0.25	0/28468	0.66	17/40222 (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
2	B	0	1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	C	N1-C2-O2	7.24	123.25	118.90
1	A	182	C	N1-C2-O2	7.10	123.16	118.90
1	A	266	U	P-O3'-C3'	6.29	127.25	119.70
1	A	367	A	P-O3'-C3'	6.00	126.90	119.70
8	H	90	LEU	CA-CB-CG	5.92	128.91	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	C	N3-C2-O2	-5.68	117.93	121.90
1	A	305	G	P-O3'-C3'	5.68	126.51	119.70
1	A	182	C	N3-C2-O2	-5.64	117.95	121.90
1	A	135	C	C2-N1-C1'	5.63	124.99	118.80
1	A	182	C	C2-N1-C1'	5.55	124.90	118.80
1	A	206	C	N1-C2-O2	5.51	122.21	118.90
1	A	329	U	C2-N1-C1'	5.31	124.07	117.70
9	J	36	LEU	CA-CB-CG	5.30	127.49	115.30
1	A	169	A	P-O3'-C3'	5.30	126.06	119.70
1	A	197	C	C5-C6-N1	5.19	123.60	121.00
1	A	271	C	C6-N1-C2	-5.13	118.25	120.30
1	A	340	C	N3-C2-O2	-5.07	118.35	121.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	363	LYS	Peptide

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	7861	0	3964	68	0
2	B	6389	0	6465	73	0
3	C	1435	0	1478	16	0
4	D	1886	0	1939	23	0
5	E	1141	0	1204	22	0
6	F	1272	0	1305	10	0
7	G	961	0	1005	10	0
8	H	1079	0	1037	9	0
9	I	1881	0	1975	24	0
9	J	2260	0	2351	29	0
10	K	1024	0	1058	19	0
11	K	1	0	0	0	0
All	All	27190	0	23781	255	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 5.

All (255) 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:A:342:C:C5	5:E:2:VAL:HA	1.96	0.99
5:E:3:ARG:H	5:E:3:ARG:HD2	1.31	0.94
1:A:191:A:O2'	1:A:192:C:OP1	1.88	0.90
1:A:151:C:O2'	1:A:190:U:OP1	1.91	0.87
1:A:193:C:H2'	1:A:194:G:C8	2.18	0.79
1:A:193:C:H2'	1:A:194:G:H8	1.50	0.75
1:A:342:C:H5	5:E:2:VAL:HA	1.51	0.73
6:F:88:ILE:O	6:F:91:MET:HB3	1.89	0.72
1:A:211:A:N6	1:A:244:G:H21	1.93	0.66
1:A:191:A:HO2'	1:A:192:C:P	2.18	0.66
1:A:211:A:N6	1:A:244:G:N2	2.46	0.64
1:A:191:A:H4'	1:A:192:C:OP2	1.97	0.64
2:B:297:ILE:HG12	2:B:367:ILE:HG22	1.79	0.63
1:A:37:U:OP2	7:G:47:ARG:NH2	2.32	0.63
2:B:441:PRO:HB3	7:G:126:LEU:HD11	1.83	0.61
1:A:343:G:OP2	5:E:3:ARG:NH2	2.34	0.61
9:J:246:THR:HG22	9:J:247:ILE:HG13	1.83	0.60
2:B:626:ARG:HH12	4:D:22:LYS:HE2	1.66	0.60
2:B:207:ASN:HD21	2:B:211:SER:HB2	1.67	0.59
2:B:646:LEU:HD23	2:B:704:ASN:HB3	1.84	0.59
5:E:29:ALA:HB1	5:E:141:ILE:HD13	1.85	0.59
5:E:29:ALA:HB2	5:E:140:ILE:HG22	1.85	0.58
2:B:454:ARG:HG3	2:B:567:PRO:HB3	1.86	0.58
1:A:191:A:O3'	1:A:192:C:H4'	2.04	0.58
4:D:227:GLY:H	4:D:251:ASP:HB3	1.69	0.58
2:B:66:ARG:NH2	5:E:81:THR:OG1	2.38	0.57
3:C:180:LYS:HG2	10:K:143:ASN:HA	1.85	0.57
9:I:67:ARG:HG3	9:I:68:ARG:HD3	1.87	0.57
10:K:115:CYS:SG	10:K:116:GLY:N	2.78	0.56
1:A:189:U:O5'	1:A:189:U:H6	1.87	0.56
1:A:98:G:H1	1:A:345:U:H3	1.51	0.56
5:E:3:ARG:HD2	5:E:3:ARG:N	2.08	0.56
3:C:44:TRP:HE1	3:C:124:LEU:HD13	1.71	0.56
3:C:126:LYS:HE2	3:C:128:ALA:HB3	1.88	0.56
2:B:762:ILE:HD13	2:B:841:PHE:HB3	1.87	0.56
4:D:198:GLY:HA3	4:D:210:MET:HB2	1.88	0.56
1:A:224:G:H2'	1:A:225:G:H8	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:A:N7	7:G:81:HIS:NE2	2.54	0.55
1:A:324:C:O2'	2:B:634:ASN:ND2	2.40	0.55
1:A:161:G:N2	4:D:18:GLU:OE2	2.40	0.54
2:B:507:GLN:NE2	2:B:508:ASN:O	2.40	0.54
2:B:297:ILE:HB	2:B:392:GLN:HB3	1.89	0.54
2:B:624:TRP:O	2:B:632:ARG:NH1	2.40	0.54
1:A:79:A:H61	6:F:134:LYS:HB3	1.72	0.54
7:G:24:LEU:HD13	7:G:29:ILE:HB	1.88	0.54
2:B:369:ARG:NH2	2:B:400:SER:OG	2.41	0.53
2:B:779:ALA:HB3	2:B:824:ILE:HB	1.90	0.53
8:H:61:ASP:HB2	8:H:64:LEU:HB3	1.90	0.53
10:K:51:ARG:NH2	10:K:57:ASP:OD1	2.41	0.53
10:K:51:ARG:HD3	10:K:58:ALA:H	1.73	0.53
1:A:134:G:H1	1:A:139:U:H3	1.54	0.53
2:B:90:GLN:HA	2:B:97:ARG:HH22	1.74	0.53
3:C:25:ASN:HD22	3:C:26:PRO:HD2	1.74	0.53
1:A:41:C:N4	7:G:18:HIS:O	2.42	0.53
6:F:37:ASN:ND2	6:F:42:THR:OG1	2.42	0.53
2:B:484:GLN:O	2:B:488:ASN:ND2	2.42	0.53
3:C:151:LYS:HA	3:C:154:VAL:HB	1.91	0.52
9:J:177:VAL:HG11	9:J:214:LEU:HD13	1.91	0.52
10:K:90:CYS:SG	10:K:91:LYS:N	2.81	0.52
2:B:88:ILE:HD12	2:B:115:ARG:HD3	1.90	0.52
9:I:120:SER:HA	9:I:124:LEU:HD12	1.90	0.52
2:B:772:ARG:HE	2:B:808:HIS:HD2	1.57	0.52
10:K:114:MET:HA	10:K:120:VAL:HG12	1.91	0.52
2:B:143:ASN:H	2:B:146:GLN:HE21	1.58	0.51
1:A:211:A:C6	1:A:244:G:N2	2.78	0.51
4:D:180:GLY:H	4:D:200:VAL:HB	1.75	0.51
5:E:2:VAL:HG13	5:E:2:VAL:O	2.10	0.51
2:B:108:ILE:O	2:B:113:ARG:NE	2.43	0.51
10:K:62:LEU:O	10:K:65:ASN:HB2	2.11	0.51
9:I:158:ILE:HB	9:I:191:ILE:HG22	1.91	0.51
9:I:219:ASP:OD2	9:I:220:ARG:NH1	2.43	0.51
2:B:662:ARG:HD3	2:B:665:GLN:HE21	1.76	0.51
1:A:208:G:N1	1:A:211:A:OP2	2.42	0.51
4:D:85:SER:HA	10:K:64:ARG:HH11	1.75	0.51
9:I:157:GLU:HA	9:I:190:VAL:O	2.10	0.50
1:A:193:C:H6	1:A:193:C:O5'	1.93	0.50
5:E:27:SER:HB3	5:E:30:ASP:HB2	1.93	0.50
1:A:54:A:H61	1:A:63:C:H42	1.58	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:45:ILE:HG23	5:E:74:LEU:HD23	1.93	0.50
5:E:116:PRO:HB2	9:I:205:LEU:HD23	1.94	0.50
1:A:211:A:H61	1:A:244:G:N2	2.09	0.49
2:B:215:THR:HG21	2:B:497:PRO:HB3	1.94	0.49
9:J:105:ASP:O	9:J:240:ASN:ND2	2.44	0.49
1:A:91:A:C2'	1:A:92:U:H5'	2.41	0.49
5:E:9:ILE:HG12	5:E:119:VAL:HG22	1.93	0.49
9:J:248:VAL:HG12	9:J:250:GLY:H	1.78	0.49
2:B:403:LEU:HB3	2:B:408:ALA:HB1	1.94	0.49
2:B:597:TYR:H	2:B:601:ASP:HB2	1.78	0.49
1:A:171:A:N6	4:D:80:ARG:O	2.45	0.49
2:B:605:THR:HG23	2:B:608:GLY:H	1.77	0.49
3:C:93:VAL:HG21	3:C:108:ILE:HG21	1.93	0.49
2:B:306:GLU:HG2	2:B:363:LYS:HE2	1.95	0.49
1:A:82:A:OP1	2:B:772:ARG:NH1	2.45	0.49
2:B:286:LEU:HB3	2:B:581:ILE:HD11	1.95	0.49
10:K:27:THR:OG1	10:K:30:ASN:ND2	2.45	0.49
2:B:157:LEU:HD21	2:B:229:LYS:HD3	1.95	0.49
8:H:39:ASN:ND2	8:H:51:GLU:OE2	2.46	0.49
9:J:48:ASN:HB2	9:J:51:GLU:HG3	1.94	0.49
4:D:179:ASN:HB3	10:K:66:TYR:HE1	1.78	0.49
10:K:110:ALA:HB1	10:K:125:ILE:H	1.77	0.49
3:C:72:PHE:HB3	10:K:105:ILE:HG21	1.95	0.48
1:A:188:U:H2'	1:A:189:U:C6	2.49	0.48
1:A:310:G:N2	1:A:344:A:N3	2.61	0.48
2:B:311:LYS:O	2:B:323:ASN:ND2	2.42	0.48
1:A:197:C:H2'	1:A:198:G:C8	2.49	0.48
8:H:85:ILE:HD11	9:J:164:LEU:HD23	1.96	0.48
1:A:363:U:O2'	4:D:222:CYS:O	2.27	0.48
2:B:470:VAL:HG12	7:G:21:LEU:HD13	1.96	0.48
2:B:816:GLU:OE1	2:B:818:HIS:NE2	2.44	0.48
3:C:99:ASP:O	3:C:133:LYS:NZ	2.39	0.48
1:A:23:G:H2'	1:A:31:A:H1'	1.96	0.47
2:B:371:HIS:O	2:B:375:TYR:N	2.43	0.47
2:B:766:SER:HB3	2:B:873:ILE:HD11	1.96	0.47
1:A:363:U:H4'	4:D:222:CYS:HB2	1.95	0.47
3:C:103:VAL:HB	10:K:125:ILE:HD12	1.94	0.47
1:A:96:G:O2'	1:A:158:U:O2	2.25	0.47
2:B:404:LYS:HD3	2:B:555:ASP:HB3	1.97	0.47
4:D:248:ILE:HG22	9:J:247:ILE:HG12	1.97	0.47
1:A:78:U:OP1	6:F:122:ARG:NH1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:J:83:ARG:HH11	9:J:108:ALA:HB2	1.79	0.47
2:B:282:SER:H	7:G:29:ILE:HD11	1.80	0.47
8:H:56:SER:OG	8:H:68:ARG:NE	2.40	0.47
9:J:87:ILE:HG12	9:J:110:LEU:HD23	1.96	0.47
1:A:270:A:O2'	2:B:577:ARG:NH2	2.43	0.47
2:B:665:GLN:HB3	2:B:728:ILE:HD11	1.97	0.47
9:I:63:ILE:O	9:I:68:ARG:NH2	2.47	0.47
9:I:177:VAL:HG11	9:I:214:LEU:HD13	1.96	0.47
2:B:333:LEU:HD12	2:B:359:PHE:HB2	1.95	0.47
2:B:780:ARG:NH2	2:B:859:ASN:O	2.47	0.47
2:B:294:GLY:O	2:B:369:ARG:HA	2.14	0.46
2:B:206:ASN:HB3	2:B:505:ASN:HA	1.97	0.46
1:A:165:A:H61	1:A:187:C:H42	1.63	0.46
7:G:90:ILE:HA	7:G:137:ARG:O	2.15	0.46
2:B:600:ASP:OD2	2:B:826:SER:OG	2.33	0.46
9:J:11:GLN:HB3	9:J:45:PHE:HE1	1.78	0.46
10:K:93:CYS:SG	10:K:95:ARG:NE	2.83	0.46
1:A:315:A:OP2	5:E:3:ARG:NH1	2.49	0.46
5:E:56:LEU:HD12	5:E:60:LEU:HD12	1.98	0.46
2:B:227:ARG:O	2:B:231:THR:OG1	2.34	0.46
9:J:3:VAL:HG22	9:J:40:HIS:HB2	1.97	0.46
1:A:320:A:O2'	9:I:220:ARG:NH2	2.48	0.45
4:D:277:ILE:HD11	9:J:234:LEU:HD22	1.97	0.45
1:A:322:G:H2'	1:A:323:A:H8	1.81	0.45
5:E:16:PRO:HA	5:E:17:PRO:HD3	1.84	0.45
2:B:856:LEU:HD13	2:B:865:TYR:HB3	1.99	0.45
3:C:30:GLU:HA	3:C:33:MET:HG2	1.99	0.45
7:G:104:ILE:HD13	7:G:122:LYS:HD2	1.98	0.45
9:I:11:GLN:HB3	9:I:45:PHE:HE1	1.80	0.45
9:I:158:ILE:HD13	9:I:177:VAL:HG13	1.98	0.45
2:B:288:MET:HB2	2:B:581:ILE:HD12	1.97	0.45
1:A:127:U:H3	1:A:146:G:H1	1.64	0.45
2:B:697:VAL:HG11	2:B:755:LEU:HD11	1.99	0.45
9:J:79:LYS:HE2	9:J:261:ASN:HA	1.99	0.45
1:A:46:G:H21	6:F:56:LYS:HA	1.82	0.45
1:A:100:A:OP2	2:B:249:ARG:NH1	2.49	0.45
10:K:63:ALA:HA	10:K:66:TYR:HD2	1.82	0.45
9:J:156:LEU:HD12	9:J:184:SER:HB2	1.98	0.45
9:I:64:ASP:HB3	9:I:67:ARG:HG2	1.98	0.44
9:I:105:ASP:OD1	9:I:240:ASN:ND2	2.50	0.44
1:A:63:C:H2'	1:A:64:G:H8	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:652:GLU:H	2:B:656:PHE:HE1	1.65	0.44
4:D:177:ASP:OD2	4:D:179:ASN:ND2	2.50	0.44
1:A:92:U:O5'	1:A:92:U:H6	1.99	0.44
2:B:110:LYS:HA	2:B:113:ARG:HD2	1.99	0.44
1:A:24:A:O3'	9:I:1:MET:N	2.51	0.44
2:B:335:GLN:HE21	2:B:552:LYS:H	1.64	0.44
1:A:193:C:O5'	1:A:193:C:C6	2.70	0.44
9:J:26:LEU:HA	9:J:29:THR:HG22	2.00	0.44
2:B:423:LYS:HB3	2:B:427:GLN:HE22	1.81	0.44
4:D:187:LYS:HB2	4:D:229:VAL:HB	1.99	0.44
2:B:115:ARG:HA	2:B:118:ARG:HG2	1.99	0.44
4:D:181:ALA:O	4:D:199:ILE:HA	2.18	0.44
10:K:61:PRO:O	10:K:65:ASN:ND2	2.51	0.44
2:B:295:THR:HA	2:B:368:ILE:O	2.18	0.43
5:E:5:LYS:HD2	5:E:90:ARG:NH1	2.34	0.43
9:I:152:ARG:HH21	9:I:154:VAL:HG21	1.83	0.43
1:A:122:C:H42	1:A:194:G:H1	1.66	0.43
9:I:128:LEU:HD23	9:I:157:GLU:HB2	1.99	0.43
9:J:9:TRP:HA	9:J:10:PRO:HD3	1.89	0.43
9:J:89:ASP:HB3	9:J:112:ILE:HD13	2.01	0.43
1:A:13:U:H2'	1:A:14:A:C8	2.54	0.43
2:B:423:LYS:O	2:B:427:GLN:NE2	2.52	0.43
9:I:127:ASP:OD1	9:I:239:ARG:NH1	2.52	0.43
9:J:29:THR:HA	9:J:198:PRO:HB3	2.00	0.43
1:A:203:U:H3	1:A:250:G:H1	1.67	0.43
9:I:9:TRP:HB3	9:I:45:PHE:HD1	1.84	0.43
9:I:86:LEU:H	9:I:109:ALA:HA	1.83	0.43
5:E:113:ILE:HG22	9:I:202:ARG:HH12	1.83	0.43
2:B:295:THR:HB	2:B:369:ARG:HB3	2.00	0.42
9:J:125:ASP:OD1	9:J:125:ASP:N	2.50	0.42
9:J:280:ASP:H	9:J:284:LEU:HD12	1.84	0.42
1:A:35:U:H2'	1:A:36:G:H8	1.84	0.42
1:A:268:U:H4'	1:A:269:G:H5'	2.00	0.42
6:F:40:ASN:HD21	6:F:150:LEU:HD23	1.84	0.42
9:J:266:ILE:HG12	9:J:273:LYS:HD2	2.01	0.42
4:D:136:TYR:HA	4:D:139:PHE:HB2	2.02	0.42
5:E:57:SER:OG	9:J:171:ARG:NH2	2.47	0.42
5:E:67:LYS:O	9:J:178:ARG:NH1	2.52	0.42
1:A:126:U:H3	1:A:147:G:H1	1.68	0.42
2:B:252:MET:HG2	2:B:261:VAL:HA	2.02	0.42
3:C:89:VAL:O	3:C:119:VAL:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:G:H3'	4:D:86:LYS:HZ1	1.85	0.42
9:I:124:LEU:HB3	9:I:126:ILE:HG23	2.01	0.42
1:A:334:G:O3'	8:H:7:ARG:NH1	2.53	0.42
6:F:51:LYS:H	6:F:87:HIS:HD2	1.66	0.42
9:I:106:ILE:HD11	9:I:233:LEU:HD23	2.02	0.42
9:J:120:SER:HA	9:J:124:LEU:HB2	2.01	0.42
1:A:96:G:O6	2:B:99:ARG:NH1	2.53	0.42
1:A:225:G:H1	1:A:232:U:H3	1.66	0.42
8:H:13:LYS:HG3	8:H:66:TYR:HE1	1.84	0.42
2:B:340:SER:HB3	2:B:353:LEU:HD13	2.01	0.42
2:B:661:TRP:HH2	2:B:841:PHE:HB2	1.85	0.42
2:B:714:GLU:HA	2:B:717:THR:HG22	2.02	0.42
2:B:767:CYS:HA	2:B:870:TRP:HA	2.01	0.41
5:E:97:ILE:HD11	5:E:119:VAL:HG21	2.00	0.41
8:H:102:GLU:HG3	9:J:205:LEU:HD21	2.02	0.41
2:B:560:LEU:HD23	2:B:565:LEU:HA	2.01	0.41
3:C:23:LEU:HB2	10:K:34:PHE:CD1	2.56	0.41
8:H:99:ILE:HG22	9:J:205:LEU:HD23	2.02	0.41
1:A:90:A:OP2	1:A:91:A:N6	2.43	0.41
3:C:24:ASP:HA	10:K:82:LEU:HD13	2.03	0.41
3:C:26:PRO:HG3	10:K:38:ASN:HD21	1.84	0.41
3:C:75:ILE:HG23	3:C:142:LEU:HD22	2.02	0.41
3:C:95:ASN:HB3	3:C:123:GLN:HE21	1.86	0.41
6:F:95:LEU:O	6:F:98:PHE:HB3	2.20	0.41
1:A:86:G:H1'	1:A:90:A:H2	1.84	0.41
1:A:342:C:H3'	5:E:3:ARG:HH21	1.85	0.41
2:B:106:ARG:O	2:B:113:ARG:NH2	2.50	0.41
2:B:217:VAL:HG22	2:B:392:GLN:HE22	1.84	0.41
4:D:122:ILE:HG22	4:D:130:TYR:HB2	2.01	0.41
8:H:94:PRO:HG2	9:J:199:LEU:HD13	2.03	0.41
2:B:445:THR:O	2:B:482:TYR:OH	2.31	0.41
1:A:240:C:H2'	1:A:241:G:C8	2.56	0.41
1:A:324:C:H2'	1:A:325:G:H8	1.85	0.41
4:D:21:GLU:HB2	4:D:24:PHE:HD2	1.85	0.41
4:D:212:VAL:HG12	4:D:221:LYS:HG3	2.02	0.41
7:G:90:ILE:HG22	7:G:138:ILE:HG12	2.02	0.41
2:B:591:TYR:OH	2:B:837:MET:O	2.39	0.41
2:B:667:LEU:HD13	2:B:820:LEU:HD21	2.03	0.41
4:D:151:ILE:HG21	4:D:234:ILE:HD12	2.01	0.41
9:I:9:TRP:HA	9:I:10:PRO:HD3	1.92	0.41
1:A:358:U:H2'	1:A:359:G:C8	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:781:ILE:HG22	2:B:821:ILE:HG13	2.03	0.41
1:A:194:G:H2'	1:A:195:A:O4'	2.21	0.40
2:B:490:VAL:HG13	2:B:544:PRO:HB3	2.03	0.40
2:B:561:PRO:HB2	2:B:564:TRP:HD1	1.86	0.40
2:B:573:ASN:HA	2:B:578:MET:HB3	2.03	0.40
6:F:116:THR:HG23	6:F:137:ILE:HD12	2.03	0.40
1:A:183:C:H2'	1:A:184:A:C8	2.56	0.40
4:D:229:VAL:HG21	4:D:276:TYR:CD1	2.57	0.40
4:D:245:ARG:HH11	9:J:251:GLY:HA3	1.86	0.40
2:B:166:MET:HG2	2:B:167:LYS:HG3	2.03	0.40
6:F:5:VAL:HG22	6:F:120:ILE:HG22	2.03	0.40
9:I:2:LEU:HD21	9:I:204:ILE:HD12	2.03	0.40
9:J:65:ILE:O	9:J:70:GLY:N	2.53	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	
2	B	774/875 (88%)	720 (93%)	54 (7%)	0	100	100
3	C	173/195 (89%)	159 (92%)	14 (8%)	0	100	100
4	D	223/279 (80%)	220 (99%)	3 (1%)	0	100	100
5	E	144/173 (83%)	140 (97%)	4 (3%)	0	100	100
6	F	155/158 (98%)	154 (99%)	1 (1%)	0	100	100
7	G	117/140 (84%)	116 (99%)	1 (1%)	0	100	100
8	H	129/133 (97%)	119 (92%)	10 (8%)	0	100	100
9	I	240/293 (82%)	225 (94%)	14 (6%)	1 (0%)	34	70
9	J	291/293 (99%)	267 (92%)	23 (8%)	1 (0%)	41	75
10	K	126/144 (88%)	111 (88%)	15 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	2372/2683 (88%)	2231 (94%)	139 (6%)	2 (0%)	54 84

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
9	I	138	PRO
9	J	138	PRO

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
2	B	708/785 (90%)	696 (98%)	12 (2%)	60 82
3	C	167/185 (90%)	163 (98%)	4 (2%)	49 75
4	D	212/261 (81%)	211 (100%)	1 (0%)	88 95
5	E	134/160 (84%)	133 (99%)	1 (1%)	84 93
6	F	148/149 (99%)	145 (98%)	3 (2%)	55 79
7	G	110/127 (87%)	109 (99%)	1 (1%)	78 91
8	H	122/123 (99%)	121 (99%)	1 (1%)	81 92
9	I	216/258 (84%)	213 (99%)	3 (1%)	67 85
9	J	258/258 (100%)	257 (100%)	1 (0%)	91 96
10	K	114/126 (90%)	112 (98%)	2 (2%)	59 81
All	All	2189/2432 (90%)	2160 (99%)	29 (1%)	70 86

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	79	ASN
2	B	99	ARG
2	B	321	ARG
2	B	443	ARG
2	B	463	ASN

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Mol	Chain	Res	Type
2	B	505	ASN
2	B	549	ARG
2	B	632	ARG
2	B	677	ASN
2	B	801	ARG
2	B	832	ASN
2	B	858	ARG
3	C	25	ASN
3	C	73	ASN
3	C	95	ASN
3	C	144	ARG
4	D	256	ARG
5	E	3	ARG
6	F	8	ASN
6	F	37	ASN
6	F	54	ASN
7	G	129	ARG
8	H	40	ASN
9	I	68	ARG
9	I	79	LYS
9	I	213	ASN
9	J	213	ASN
10	K	38	ASN
10	K	55	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	79	ASN
2	B	146	GLN
2	B	207	ASN
2	B	218	ASN
2	B	243	HIS
2	B	246	ASN
2	B	251	HIS
2	B	335	GLN
2	B	392	GLN
2	B	427	GLN
2	B	505	ASN
2	B	590	GLN
2	B	593	ASN
2	B	595	GLN

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Mol	Chain	Res	Type
2	B	634	ASN
2	B	642	HIS
2	B	665	GLN
2	B	677	ASN
2	B	808	HIS
2	B	832	ASN
3	C	25	ASN
3	C	32	HIS
3	C	73	ASN
3	C	95	ASN
3	C	160	ASN
4	D	5	GLN
5	E	59	ASN
6	F	8	ASN
6	F	37	ASN
6	F	54	ASN
6	F	76	GLN
7	G	85	GLN
8	H	27	HIS
8	H	39	ASN
8	H	40	ASN
9	I	151	ASN
9	I	213	ASN
9	I	240	ASN
10	K	30	ASN
10	K	38	ASN
10	K	47	GLN
10	K	140	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	368/369 (99%)	100 (27%)	9 (2%)

All (100) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	16	U
1	A	17	U
1	A	24	A
1	A	26	U

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Mol	Chain	Res	Type
1	A	28	A
1	A	29	G
1	A	30	A
1	A	32	A
1	A	37	U
1	A	38	U
1	A	39	U
1	A	40	A
1	A	42	A
1	A	43	G
1	A	44	A
1	A	46	G
1	A	47	G
1	A	61	G
1	A	68	U
1	A	74	A
1	A	76	A
1	A	78	U
1	A	79	A
1	A	80	U
1	A	88	G
1	A	89	A
1	A	90	A
1	A	91	A
1	A	93	U
1	A	94	C
1	A	99	G
1	A	100	A
1	A	101	A
1	A	102	C
1	A	105	A
1	A	106	G
1	A	115	U
1	A	116	G
1	A	118	C
1	A	119	C
1	A	120	U
1	A	121	C
1	A	135	C
1	A	137	U
1	A	138	U
1	A	152	U

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Mol	Chain	Res	Type
1	A	153	G
1	A	167	A
1	A	168	A
1	A	170	U
1	A	171	A
1	A	172	U
1	A	173	A
1	A	177	G
1	A	179	G
1	A	180	A
1	A	182	C
1	A	187	C
1	A	189	U
1	A	190	U
1	A	191	A
1	A	192	C
1	A	196	C
1	A	197	C
1	A	203	U
1	A	213	U
1	A	214	C
1	A	229	A
1	A	230	G
1	A	231	G
1	A	243	U
1	A	247	A
1	A	257	G
1	A	258	U
1	A	259	U
1	A	260	G
1	A	267	U
1	A	268	U
1	A	279	C
1	A	284	A
1	A	287	G
1	A	291	G
1	A	292	A
1	A	301	G
1	A	304	U
1	A	305	G
1	A	306	U
1	A	307	U

Continued on next page...

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Mol	Chain	Res	Type
1	A	309	A
1	A	313	C
1	A	317	C
1	A	329	U
1	A	331	A
1	A	332	G
1	A	333	U
1	A	338	A
1	A	343	G
1	A	344	A
1	A	367	A
1	A	368	G

All (9) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	169	A
1	A	179	G
1	A	191	A
1	A	195	A
1	A	228	U
1	A	266	U
1	A	303	C
1	A	305	G
1	A	367	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

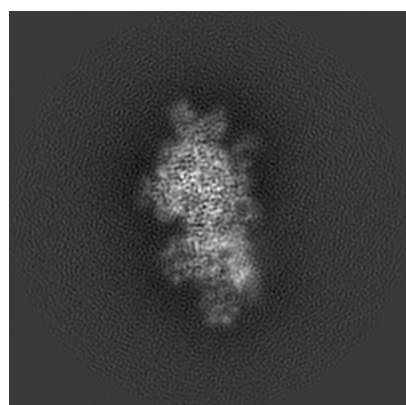
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9616. These allow visual inspection of the internal detail of the map and identification of artifacts.

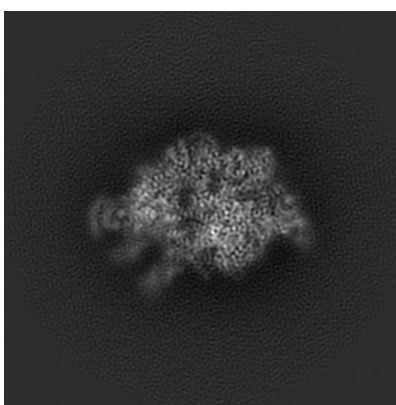
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

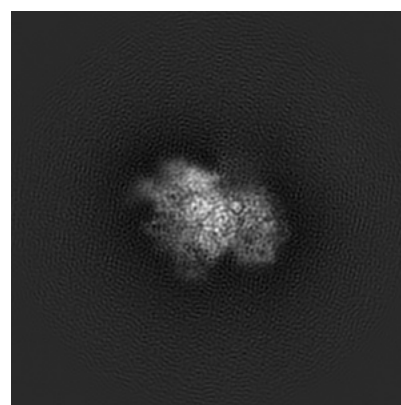
6.1.1 Primary map



X



Y

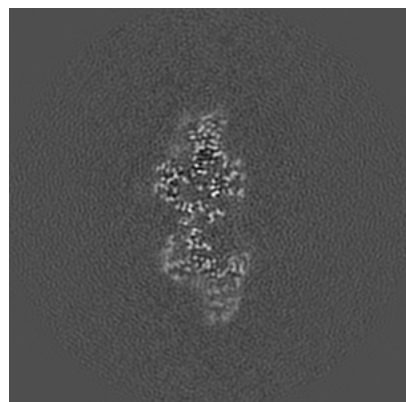


Z

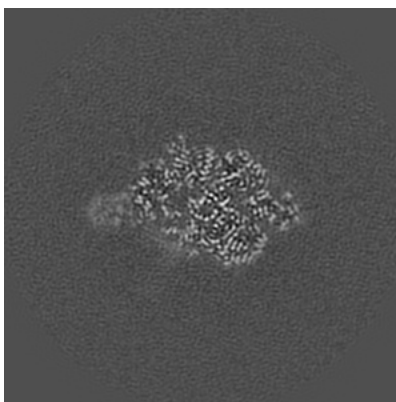
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

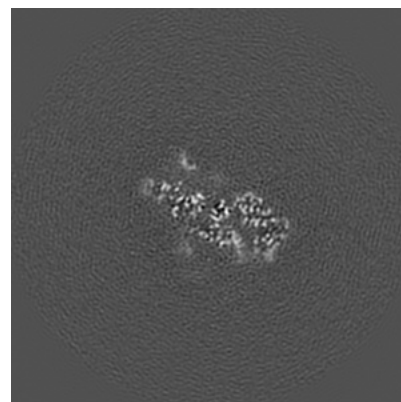
6.2.1 Primary map



X Index: 128



Y Index: 128

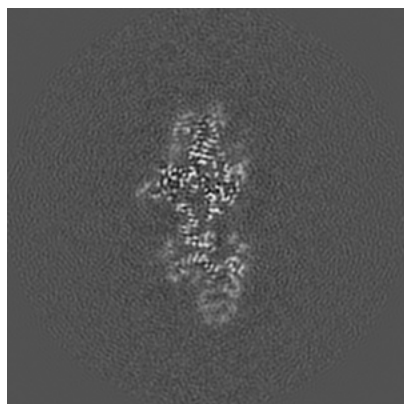


Z Index: 128

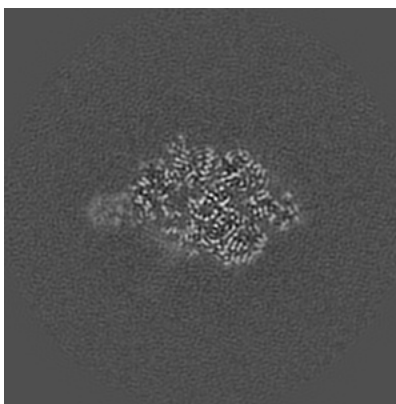
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

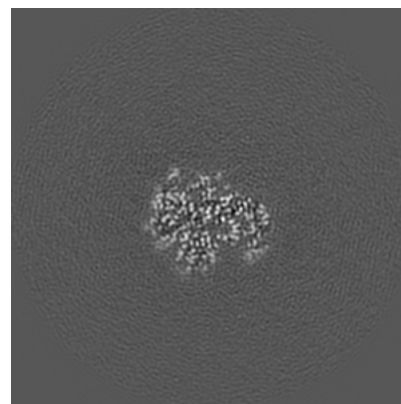
6.3.1 Primary map



X Index: 124



Y Index: 128



Z Index: 142

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

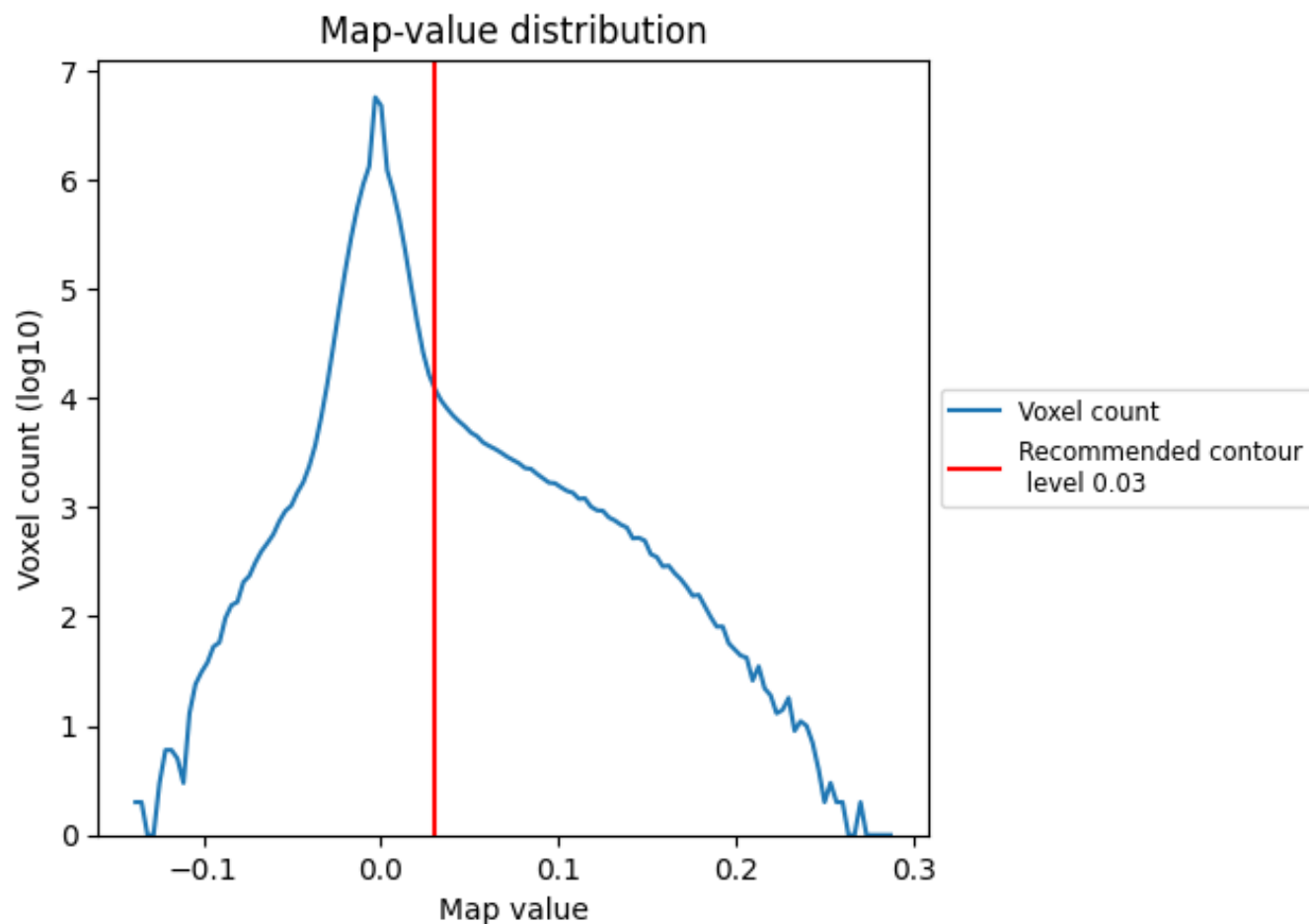
6.5 Mask visualisation

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

7 Map analysis [i](#)

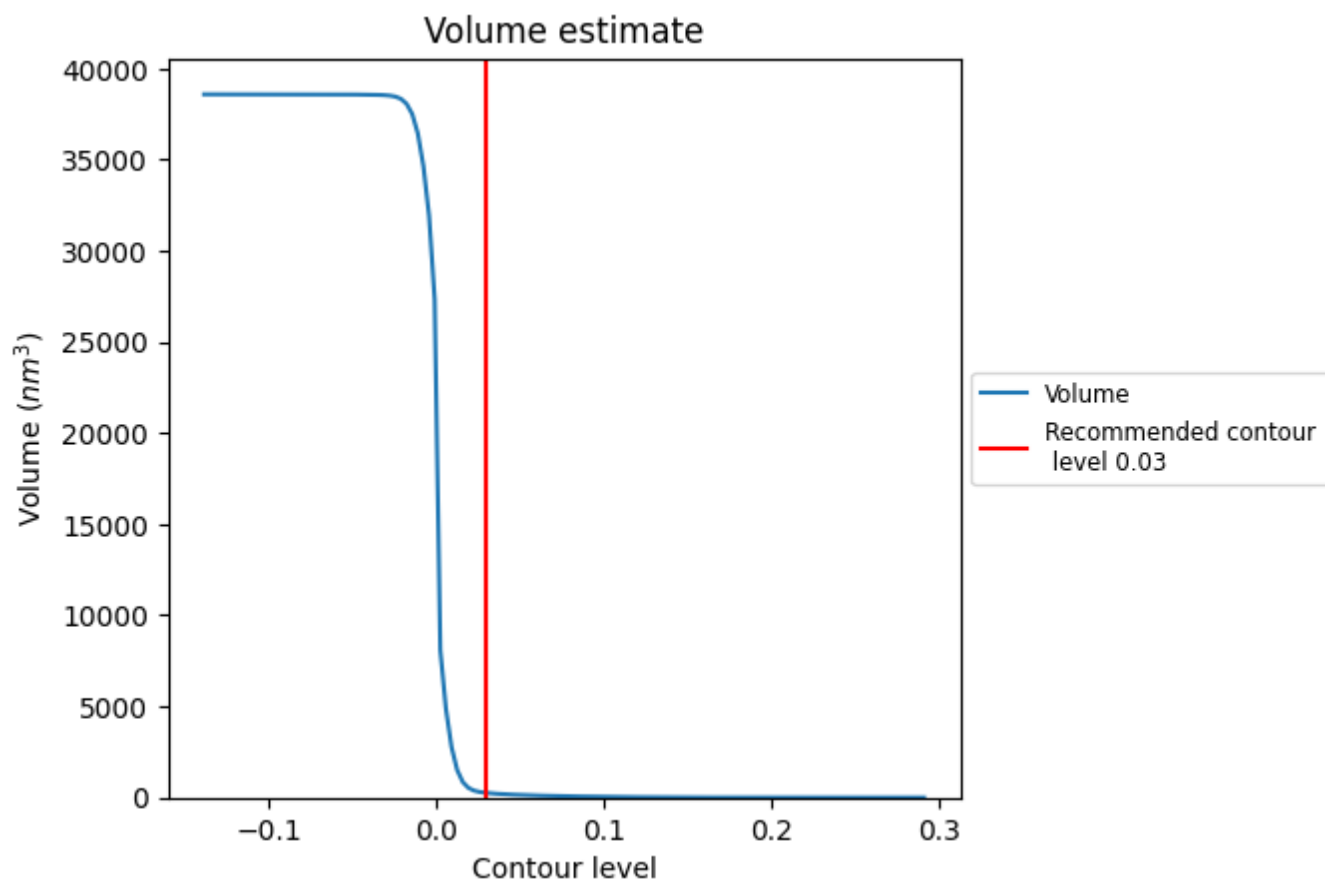
This section contains the results of statistical analysis of the map.

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.

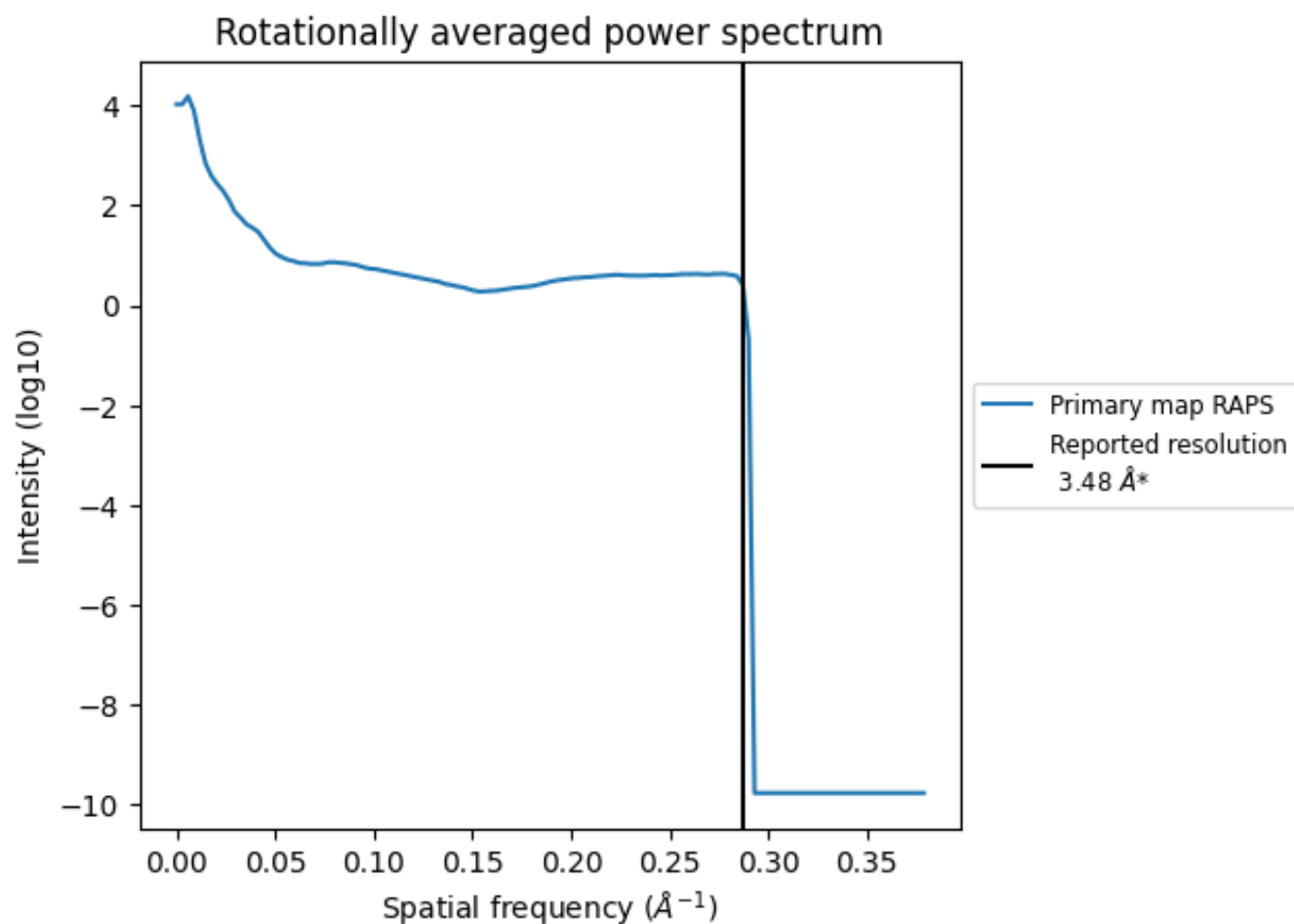
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 257 nm^3 ; this corresponds to an approximate mass of 232 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 0.287 Å⁻¹

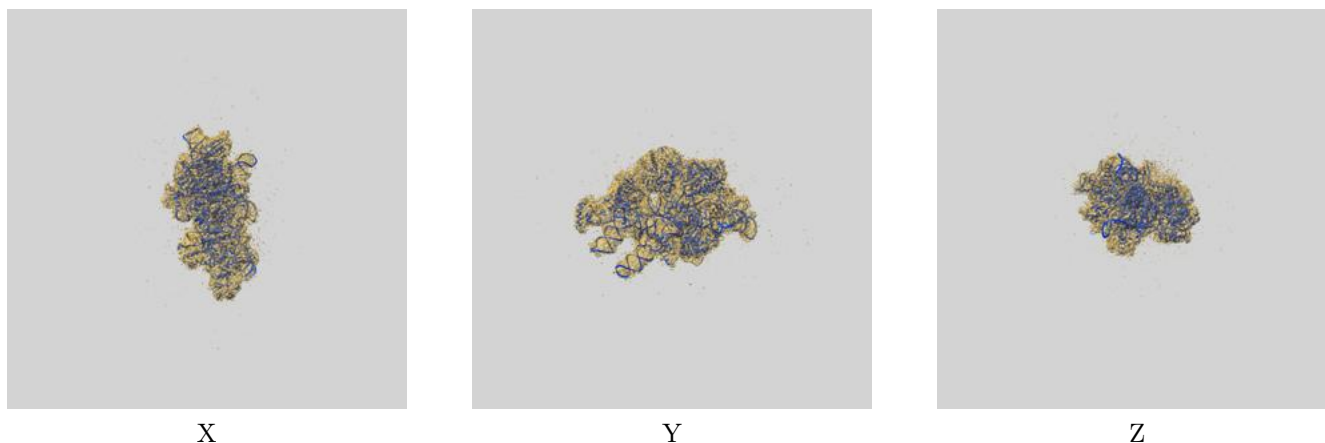
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

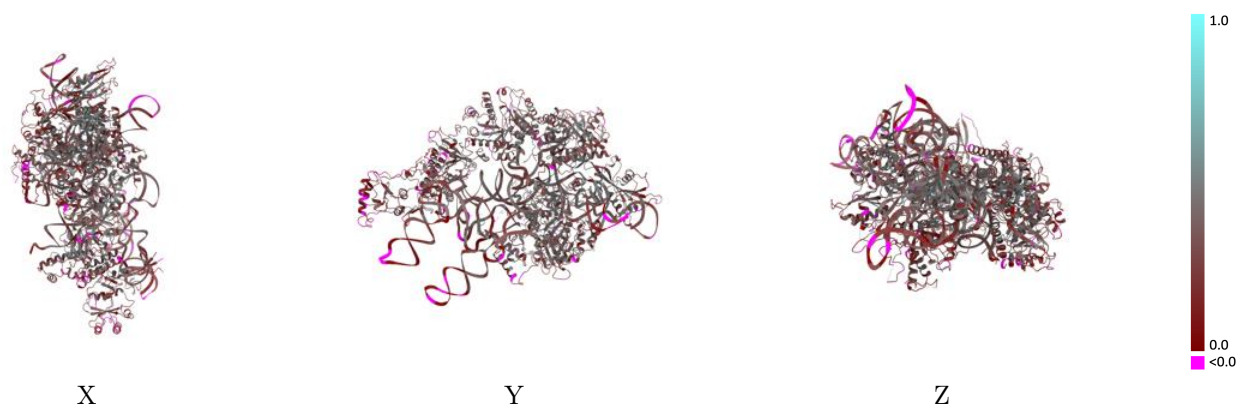
This section contains information regarding the fit between EMDB map EMD-9616 and PDB model 6AGB. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



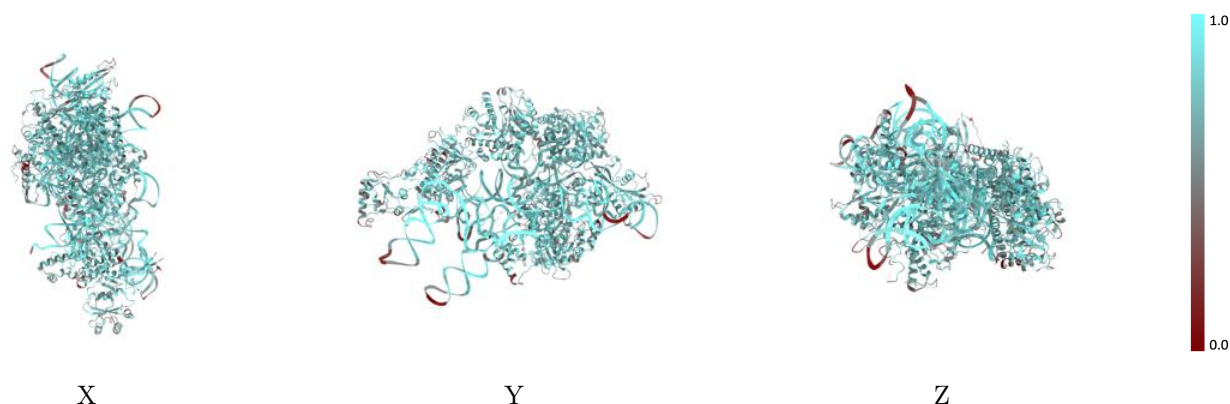
The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



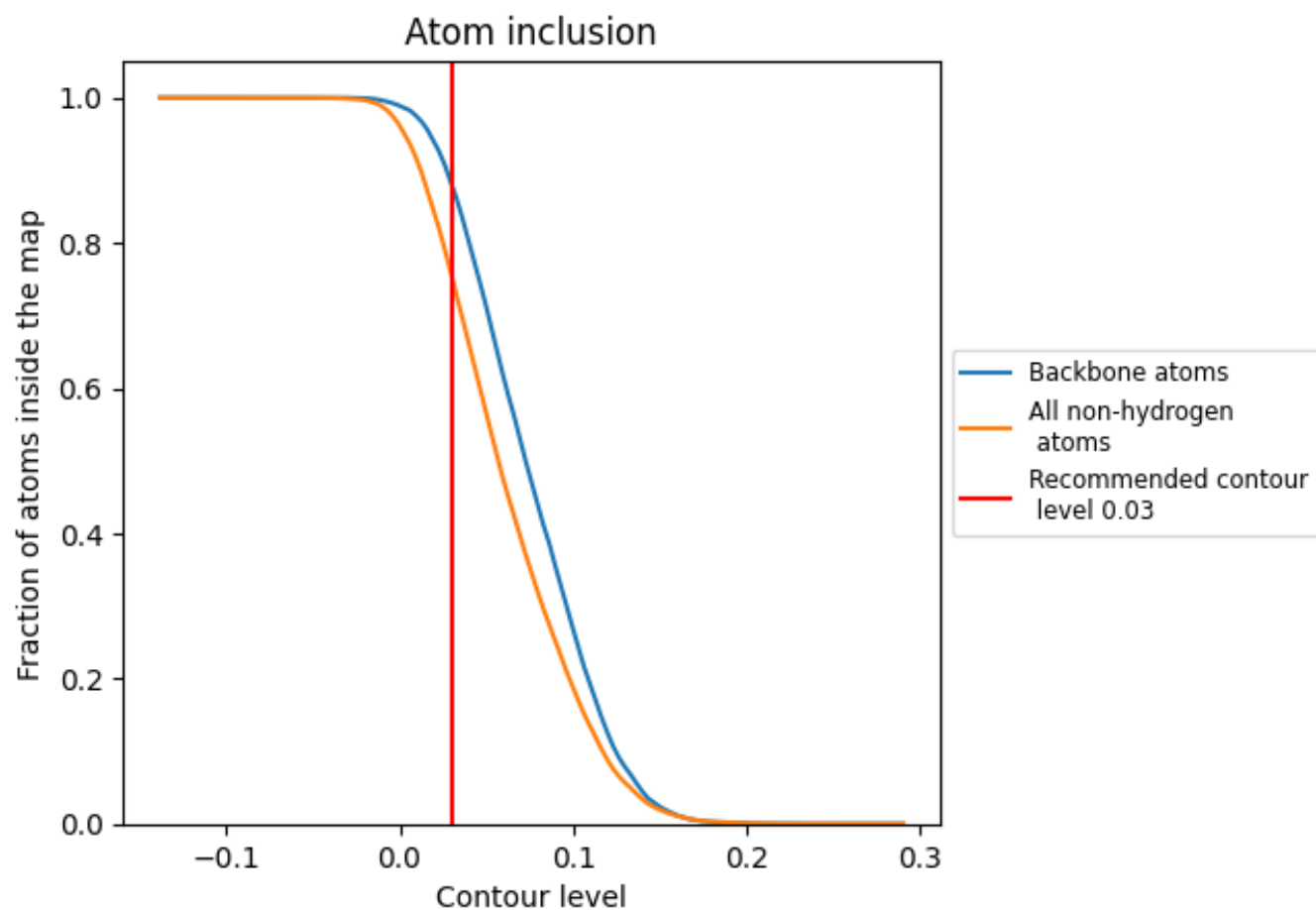
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.03).

























9.4 Atom inclusion [i](#)



At the recommended contour level, 88% 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 (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7549	 0.3110
A	 0.8144	 0.2930
B	 0.7414	 0.3230
C	 0.6416	 0.2190
D	 0.7015	 0.3050
E	 0.7741	 0.3560
F	 0.7526	 0.3370
G	 0.7840	 0.3930
H	 0.6531	 0.2510
I	 0.7391	 0.3320
J	 0.7517	 0.3480
K	 0.7269	 0.3020

