



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 03:11 AM UTC

PDB ID : 8C00 / pdb_00008c00
EMDB ID : EMD-16347
Title : Enp1TAP-S21_A population of yeast small ribosomal subunit precursors depleted of rpS21/eS21
Authors : Milkereit, P.; Poell, G.
Deposited on : 2022-12-15
Resolution : 2.90 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

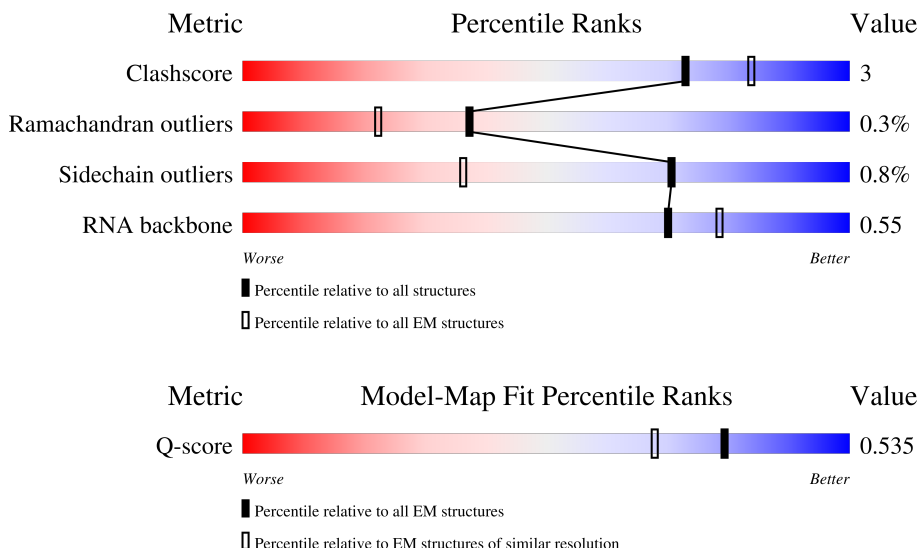
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 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.



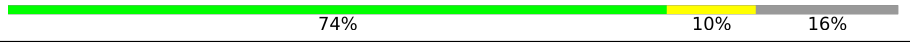
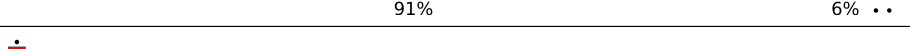
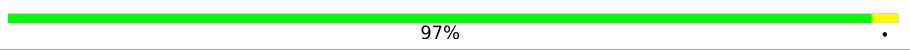
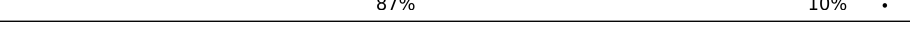
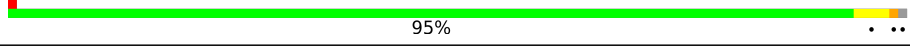
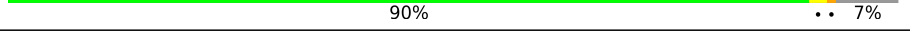
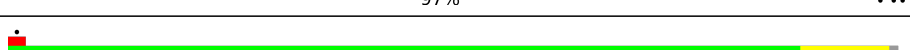
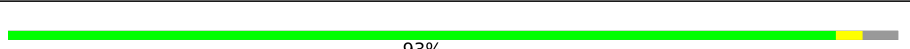
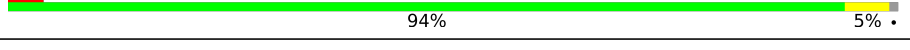
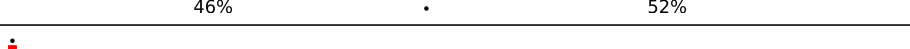
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13054 (2.40 - 3.40)

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	2	1800	
2	B	225	
3	E	142	

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Mol	Chain	Length	Quality of chain
4	F	143	
5	H	146	
6	I	144	
7	K	108	
8	L	67	
9	Q	255	
10	S	261	
11	T	236	
12	U	190	
13	V	200	
14	W	197	
15	X	156	
16	Y	151	
17	Z	137	
18	b	130	
19	c	145	
20	d	135	
21	e	483	
22	f	82	
23	g	63	
24	p	274	
25	r	425	
26	t	788	

2 Entry composition

There are 26 unique types of molecules in this entry. The entry contains 63117 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 18S rRNA precursor.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	2	1425	Total	C	N	O	P	0	0
			30408	13596	5424	9963	1425		

- Molecule 2 is a protein called rpS5.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	201	Total	C	N	O	S	0	0
			1588	996	295	294	3		

- Molecule 3 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	114	Total	C	N	O	S	0	0
			902	575	167	153	7		

- Molecule 4 is a protein called 40S ribosomal protein S16-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	F	125	Total	C	N	O	0	0
			973	625	174	174		

- Molecule 5 is a protein called 40S ribosomal protein S18-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	122	Total	C	N	O	S	0	0
			1009	636	193	178	2		

- Molecule 6 is a protein called 40S ribosomal protein S19-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	I	141	Total	C	N	O	S	0	0
			1096	684	206	204	2		

- Molecule 7 is a protein called 40S ribosomal protein S25-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	K	69	Total	C	N	O	0	0
			556	356	103	97		

- Molecule 8 is a protein called 40S ribosomal protein S28-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	L	60	Total	C	N	O	S	0	0
			471	290	93	87	1		

- Molecule 9 is a protein called 40S ribosomal protein S1-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	Q	214	Total	C	N	O	S	0	0
			1709	1084	310	311	4		

- Molecule 10 is a protein called 40S ribosomal protein S4-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	S	260	Total	C	N	O	S	0	0
			2068	1316	389	360	3		

- Molecule 11 is a protein called 40S ribosomal protein S6-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	T	226	Total	C	N	O	S	0	0
			1813	1137	350	323	3		

- Molecule 12 is a protein called 40S ribosomal protein S7-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
12	U	184	Total	C	N	O	0	0
			1481	951	265	265		

- Molecule 13 is a protein called 40S ribosomal protein S8-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	V	185	Total	C	N	O	S	0	0
			1462	907	292	261	2		

- Molecule 14 is a protein called 40S ribosomal protein S9-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	W	185	Total	C	N	O	S	0	0
			1494	943	289	261	1		

- Molecule 15 is a protein called 40S ribosomal protein S11-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	X	141	Total	C	N	O	S	0	0
			1137	730	216	188	3		

- Molecule 16 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	Y	150	Total	C	N	O	S	0	0
			1192	759	224	207	2		

- Molecule 17 is a protein called 40S ribosomal protein S14-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Z	127	Total	C	N	O	S	0	0
			926	569	185	169	3		

- Molecule 18 is a protein called 40S ribosomal protein S22-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	b	129	Total	C	N	O	S	0	0
			1021	650	188	180	3		

- Molecule 19 is a protein called 40S ribosomal protein S23-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	c	144	Total	C	N	O	S	0	0
			1121	708	220	191	2		

- Molecule 20 is a protein called 40S ribosomal protein S24-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
20	d	130	Total	C	N	O	0	0
			1046	662	203	181		

- Molecule 21 is a protein called Essential nuclear protein 1.

Mol	Chain	Residues	Atoms				AltConf	Trace
21	e	179	Total	C	N	O	0	0
			796	431	183	182		

- Molecule 22 is a protein called 40S ribosomal protein S27-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
22	f	81	Total	C	N	O	S	0
			610	382	110	113	5	0

- Molecule 23 is a protein called 40S ribosomal protein S30-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
23	g	30	Total	C	N	O	S	0
			257	164	55	37	1	0

- Molecule 24 is a protein called Pre-rRNA-processing protein PNO1.

Mol	Chain	Residues	Atoms				AltConf	Trace
24	p	180	Total	C	N	O	S	0
			1425	911	257	253	4	0

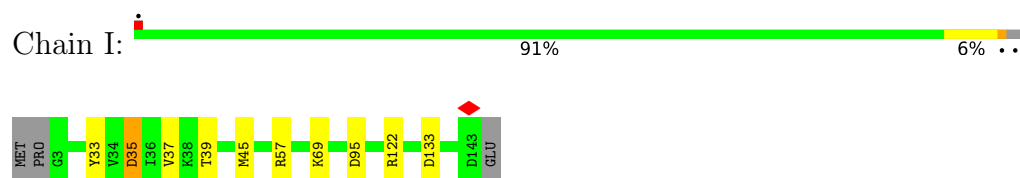
- Molecule 25 is a protein called Serine/threonine-protein kinase RIO2.

Mol	Chain	Residues	Atoms				AltConf	Trace
25	r	190	Total	C	N	O	S	0
			1574	1014	268	278	14	0

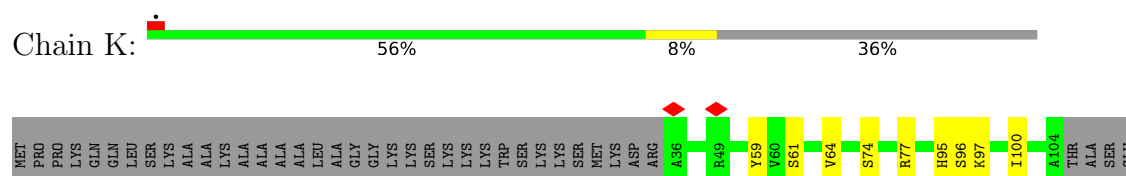
- Molecule 26 is a protein called Ribosome biogenesis protein TSR1.

Mol	Chain	Residues	Atoms				AltConf	Trace
26	t	619	Total	C	N	O	S	0
			4982	3199	873	897	13	0

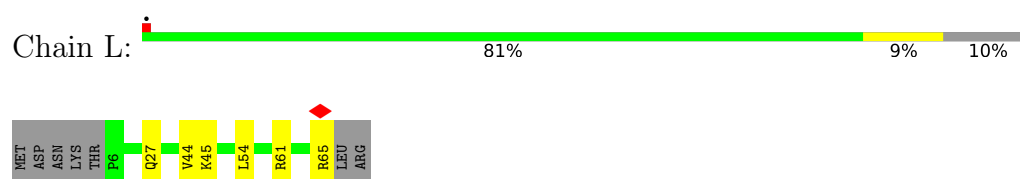
- Molecule 6: 40S ribosomal protein S19-A



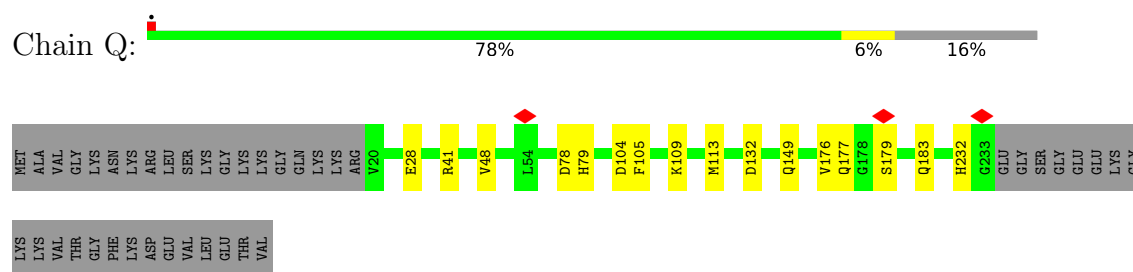
- Molecule 7: 40S ribosomal protein S25-A



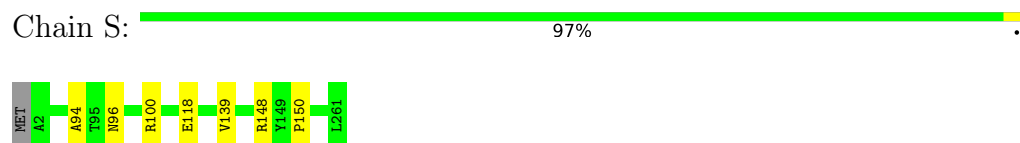
- Molecule 8: 40S ribosomal protein S28-A



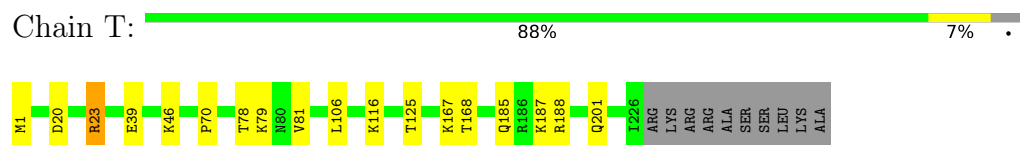
- Molecule 9: 40S ribosomal protein S1-A



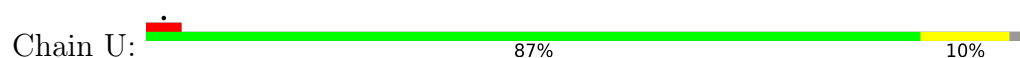
- Molecule 10: 40S ribosomal protein S4-A



- Molecule 11: 40S ribosomal protein S6-A



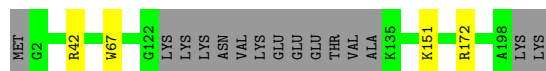
- Molecule 12: 40S ribosomal protein S7-A





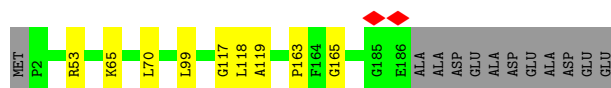
- Molecule 13: 40S ribosomal protein S8-A

Chain V: 90% 8%



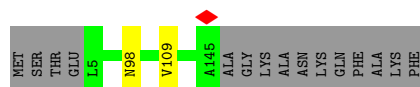
- Molecule 14: 40S ribosomal protein S9-A

Chain W: 89% 5% 6%



- Molecule 15: 40S ribosomal protein S11-A

Chain X: 89% 10%



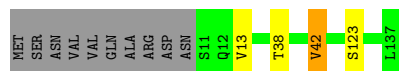
- Molecule 16: 40S ribosomal protein S13

Chain Y: 95%



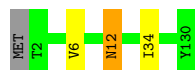
- Molecule 17: 40S ribosomal protein S14-A

Chain Z: 90% 7%



- Molecule 18: 40S ribosomal protein S22-A

Chain b: 97%

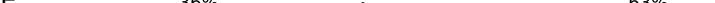


- Molecule 19: 40S ribosomal protein S23-A

Amino Acid	Count
MET	1
G2	1
N27	1
S41	1
P42	1
I59	1
E60	1
Q63	1
P64	1
M65	1
I68	1
R69	1
K70	1
R73	1
N79	1
K82	1
D90	1
G91	1
K110	1
D116	1
S145	1

- Chain d: 93% . .

MET
SER
D3
H29
T121
R124 L125
R132
ASN
ALA
ASP

- Chain e:  36% 63%

MET	ARG	ALA	ALA	ARG	SER	SER	THR	LYS	ALA	ALA	GLN	LYS	GLN	GLN	ASP	PRO	ASP	ASP	LEU	LEU	LEU	ALA	ALA	GLN	GLY	THR	LYS	LYS	LYS	LYS	LEU	LEU	GLN	ASN	ASN	LYS	LYS	LYS	LYS	LEU	LEU	GLN	ASP	ASN	ASN	ALA	ALA	ASN	GLU	GLU	ASP	GLY	TYR	ILE	ILE	ASP	SER	LYS	ALA	ALA	SER	PER
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LYS ILE LEU LEU GLN GLN ALA LYS LYS GLU GLU GLN GLN ASP ASP LEU ILE LEU GLU GLY GLY GLU GLU LEU LEU ALA GLU GLU SER SER PHE ARG ASN LYS LYS GLN GLN PHE PHE GLU GLU ALA ALA ARG ARG PHE PHE THR THR MET MET SER SER TYR TYR ASP ASP ASP ASP ASP ASP ASP ASP ASP ASP GLU GLU GLU GLU GLU GLU GLU GLU GLU GLU ALA ALA PHE PHE GLY GLY GLU GLU ASP ASP ILE ILE SER SER ASP ASP PHE PHE GLU GLU PRO PRO

GLY	ASP	TYR	LYS	GLU	GLU	GLU	ILE	VAL	GLU	ILE	ASP	GLU	GLY	ASP	ALA	ALA	MET	PHE	GLU	GLN	TYR	PHE	LYS	LYS	SER	SER	ASP	ASP	PHE	ASN	SER	LEU	LEU	ALA	ALA	ASP	LYS	ILE	ILE	ARG	GLU	LYS	GLU	SER	GLN	VAL	GLU	ASP	MET	GLN	ASP	SER
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GLU	PRO	LEU	ALA	ASN	GLU	GLN	ASN	THR	SER	ARG	GLY	ASN	ILE	SER	SER	GLY	LEU	LYS	SER	GLY	GLU	GLY	VAL	ALA	LEU	PRO	GLU	GLU	LYS	VAL	ILE	LYS	A213	T216	T223	L232	T235	T237	S238	V248	V258	A262	V267	SER	ASN	LEU	THR	ALA	K273	T282
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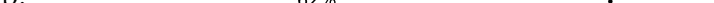
ASP	GLN	ARG	ASP	PHE	LEU	LEU	GLU	THR	VAL	ARG	GLN	ARG	GLY	HIS	LYS	ASP	ILE	GLY	PRO	ILE	ARG	ARG	GLU	LEU	LEU	ALA	ALA	GLY	ALA	SER	ARG	GLU	PHE	VAL	ASP	PRO	GLN	GLN	GLU	ALA	ASN	ASP	ASP	LEU	MET	ILE	ILE	ASP	VAL	VAL	ASN
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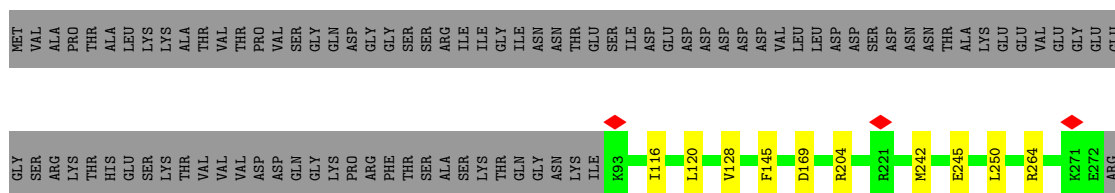
- Chain f: 94% 5%

- Chain g: 46% 52%

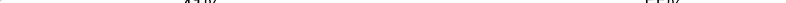
MET	ALA	VAL	HIS	GLY	SER	LEU	ALA	ARG	ALA	GLY	LYS	VAL	LYS	SER	GLN	THR	PRO	LYS	VAL	GLU	LYS	THR	E25	K31	V47	THR	LEU	VAL	ASN	GLY	LYS	R54	P60	SER	VAL	GLN
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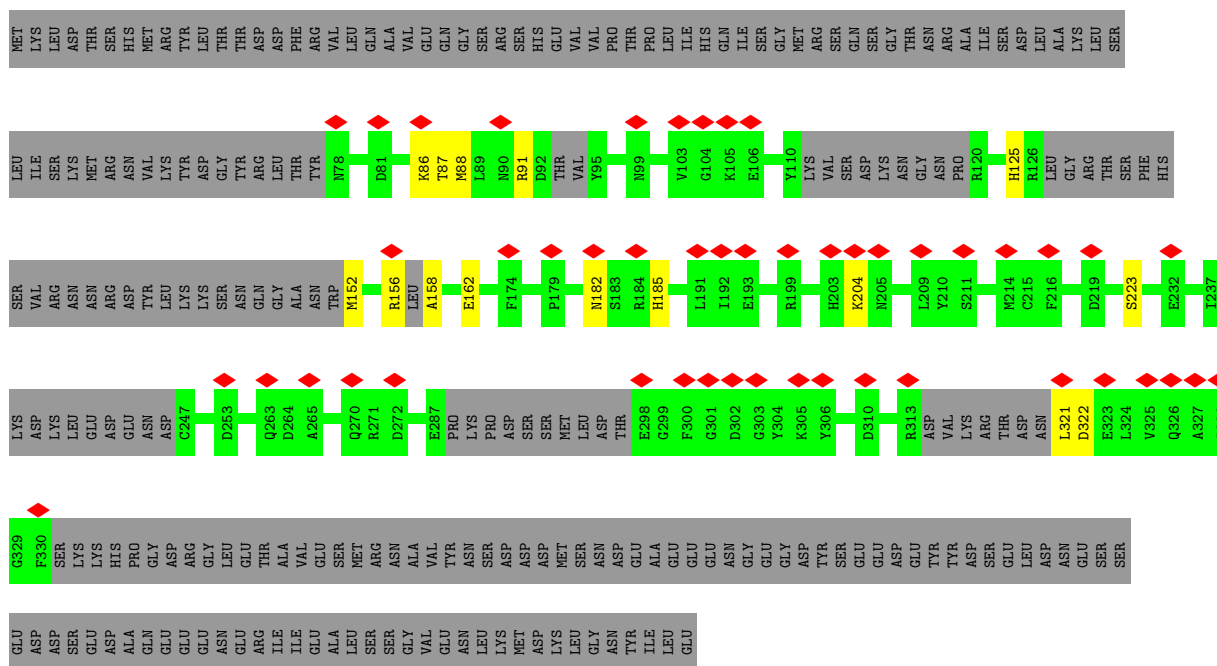
- WORLDWIDE
PDB
PROTEIN DATA BANK

Chain p:  62% 34%

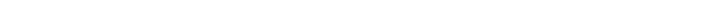


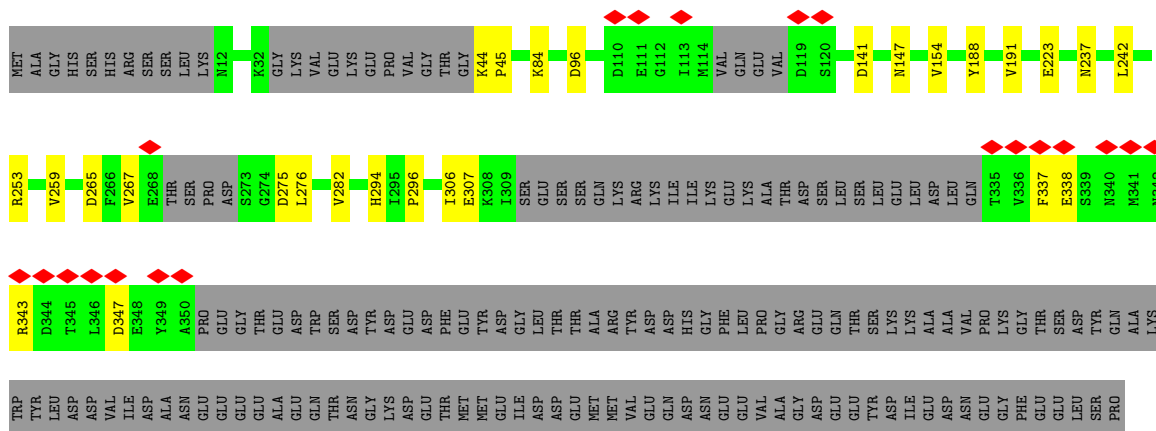
- Molecule 25: Serine/threonine-protein kinase RIO2

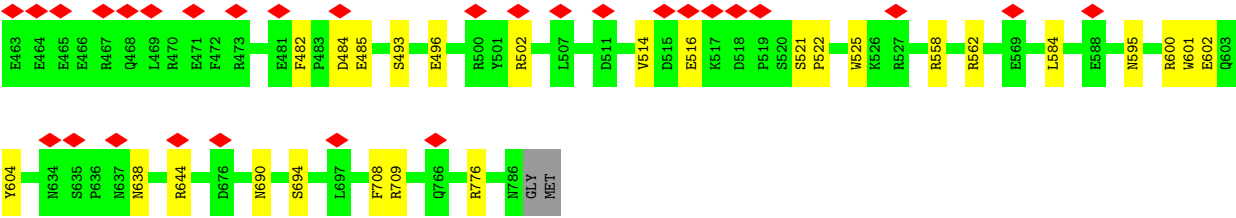
Chain r: 



- Molecule 26: Ribosome biogenesis protein TSR1

Chain t: 





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	132563	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	JEOL CRYO ARM 200	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	50000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.048	Depositor
Minimum map value	-0.010	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.007	Depositor
Map size (Å)	387.2, 387.2, 387.2	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.96800005, 0.96800005, 0.96800005	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	2	0.38	0/33998	0.53	7/52913 (0.0%)
2	B	0.19	0/1607	0.39	0/2172
3	E	0.18	0/921	0.41	0/1236
4	F	0.18	0/990	0.41	0/1335
5	H	0.19	0/1027	0.41	0/1383
6	I	0.20	0/1113	0.41	0/1494
7	K	0.15	0/564	0.36	0/758
8	L	0.19	0/473	0.45	0/634
9	Q	0.27	0/1735	0.47	0/2335
10	S	0.35	0/2109	0.50	0/2839
11	T	0.28	0/1837	0.43	0/2455
12	U	0.24	0/1506	0.46	0/2028
13	V	0.35	0/1487	0.49	0/1988
14	W	0.35	0/1519	0.47	0/2035
15	X	0.36	0/1163	0.43	0/1568
16	Y	0.33	0/1215	0.43	0/1638
17	Z	0.25	0/937	0.48	0/1261
18	b	0.37	0/1038	0.43	0/1395
19	c	0.32	0/1139	0.46	0/1518
20	d	0.33	0/1060	0.46	0/1412
21	e	0.10	0/807	0.34	2/1025 (0.2%)
22	f	0.29	0/620	0.41	0/838
23	g	0.34	0/262	0.50	0/346
24	p	0.26	0/1451	0.43	0/1955
25	r	0.12	0/1605	0.36	0/2141
26	t	0.17	0/5092	0.39	0/6868
All	All	0.33	0/67275	0.49	9/97570 (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
6	I	0	1
9	Q	0	1
10	S	0	1
16	Y	0	1
26	t	0	1
All	All	0	5

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	143	G	N9-C1'-C2'	6.29	121.43	112.00
1	2	514	G	P-O3'-C3'	5.95	129.12	120.20
1	2	9	U	C5'-C4'-O4'	5.33	117.09	109.10
1	2	514	G	C4'-C3'-O3'	5.32	117.37	109.40
1	2	1653	C	N1-C1'-C2'	5.18	119.78	112.00
1	2	9	U	C1'-O4'-C4'	-5.17	104.53	109.70
1	2	1492	A	C2'-C3'-O3'	5.13	117.19	109.50
21	e	235	VAL	CA-C-N	5.11	123.46	120.24
21	e	235	VAL	C-N-CA	5.11	123.46	120.24

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
6	I	57	ARG	Sidechain
9	Q	132	ASP	Peptide
10	S	148	ARG	Sidechain
16	Y	3	ARG	Sidechain
26	t	253	ARG	Sidechain

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	2	30408	0	15313	251	0
2	B	1588	0	1657	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	902	0	934	6	0
4	F	973	0	1029	3	0
5	H	1009	0	1029	10	0
6	I	1096	0	1110	5	0
7	K	556	0	596	5	0
8	L	471	0	505	4	0
9	Q	1709	0	1784	8	0
10	S	2068	0	2154	3	0
11	T	1813	0	1905	12	0
12	U	1481	0	1572	10	0
13	V	1462	0	1486	3	0
14	W	1494	0	1573	5	0
15	X	1137	0	1207	0	0
16	Y	1192	0	1255	4	0
17	Z	926	0	950	2	0
18	b	1021	0	1060	2	0
19	c	1121	0	1196	9	0
20	d	1046	0	1112	2	0
21	e	796	0	296	1	0
22	f	610	0	633	1	0
23	g	257	0	281	2	0
24	p	1425	0	1502	6	0
25	r	1574	0	1545	8	0
26	t	4982	0	4954	26	0
All	All	63117	0	48638	361	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 3.

All (361) 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)
1:2:1690:G:H21	1:2:1712:A:N6	1.52	1.05
1:2:1690:G:N2	1:2:1712:A:H61	1.57	1.02
1:2:1588:G:H1	1:2:1608:U:H3	1.03	0.94
1:2:1690:G:H21	1:2:1712:A:H61	0.90	0.89
26:t:521:SER:OG	26:t:525:TRP:NE1	2.10	0.84
1:2:1081:A:N3	1:2:1082:C:N4	2.28	0.80
1:2:65:A:H2	1:2:84:A:H62	1.30	0.79
1:2:567:A:H62	1:2:576:G:N2	1.79	0.79
1:2:901:G:H2'	1:2:902:G:C8	2.20	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1748:G:H2'	1:2:1749:A:C8	2.20	0.76
1:2:567:A:H62	1:2:576:G:H21	1.33	0.74
1:2:1690:G:N3	1:2:1712:A:C6	2.54	0.74
1:2:1690:G:N2	1:2:1712:A:N6	2.24	0.73
26:t:141:ASP:OD1	26:t:147:ASN:ND2	2.23	0.71
1:2:1524:A:H2'	1:2:1525:A:C8	2.26	0.70
1:2:1499:G:OP1	6:I:122:ARG:NH1	2.24	0.70
26:t:307:GLU:OE2	26:t:558:ARG:NH1	2.24	0.70
1:2:1206:U:OP2	1:2:1207:C:O2'	2.09	0.70
26:t:502:ARG:NH2	26:t:516:GLU:OE2	2.24	0.70
1:2:1240:U:O2'	1:2:1242:A:N6	2.25	0.69
1:2:846:G:H2'	1:2:847:A:H8	1.58	0.69
1:2:895:G:H1	1:2:917:U:H3	1.41	0.68
1:2:992:A:H2	1:2:1012:U:H3	1.41	0.68
1:2:821:U:H3	1:2:852:C:H42	1.42	0.68
4:F:50:GLU:OE1	4:F:82:ARG:NH2	2.26	0.68
1:2:846:G:H2'	1:2:847:A:C8	2.29	0.67
1:2:1672:G:H2'	1:2:1673:G:C8	2.30	0.67
1:2:1690:G:N3	1:2:1712:A:N1	2.43	0.67
1:2:1160:A:H2'	1:2:1161:C:H6	1.59	0.66
12:U:7:LYS:O	12:U:42:GLN:NE2	2.28	0.65
19:c:42:PRO:O	19:c:79:ASN:ND2	2.30	0.65
3:E:18:ARG:NH1	5:H:90:ASN:OD1	2.30	0.65
1:2:1592:A:H2'	1:2:1593:A:H8	1.61	0.64
1:2:1653:C:H3'	1:2:1654:G:H8	1.60	0.64
1:2:567:A:N6	1:2:576:G:H21	1.96	0.64
1:2:1207:C:H5'	1:2:1208:A:C8	2.32	0.63
1:2:1525:A:H2'	1:2:1526:A:C8	2.33	0.63
1:2:554:C:H1'	1:2:555:A:H2	1.63	0.63
1:2:1653:C:H3'	1:2:1654:G:C8	2.34	0.63
1:2:513:U:H3	1:2:538:A:H2	1.43	0.63
1:2:1690:G:C2	1:2:1712:A:N1	2.67	0.63
25:r:156:ARG:O	25:r:158:ALA:N	2.32	0.62
1:2:828:U:H2'	1:2:829:A:H8	1.64	0.62
1:2:1727:G:H2'	1:2:1728:A:C8	2.34	0.62
1:2:1081:A:O2'	1:2:1083:G:N7	2.32	0.62
6:I:37:VAL:HG12	6:I:39:THR:H	1.64	0.62
1:2:851:U:H2'	1:2:852:C:C6	2.35	0.62
1:2:924:A:H2'	1:2:925:G:C8	2.35	0.62
1:2:1752:U:H2'	1:2:1753:A:C8	2.36	0.61
1:2:545:A:H2'	23:g:31:LYS:HD3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:94:THR:HG22	2:B:114:ILE:HG13	1.84	0.60
1:2:284:G:H22	11:T:188:ARG:NH1	1.99	0.60
1:2:1368:G:H5''	6:I:69:LYS:HE2	1.81	0.60
1:2:1592:A:H2'	1:2:1593:A:C8	2.37	0.60
1:2:1555:A:OP2	3:E:47:ARG:NH1	2.35	0.59
1:2:1501:C:N4	1:2:1502:G:O6	2.35	0.59
1:2:1690:G:H1'	1:2:1711:C:H5''	1.85	0.59
26:t:306:ILE:O	26:t:337:PHE:N	2.28	0.59
1:2:1002:G:N2	1:2:1005:A:O2'	2.35	0.59
1:2:1649:G:OP1	19:c:82:LYS:NZ	2.35	0.58
1:2:563:U:H2'	1:2:564:G:C8	2.37	0.58
1:2:1254:U:H2'	1:2:1255:G:C8	2.38	0.58
26:t:343:ARG:HH22	26:t:562:ARG:NH2	2.01	0.58
1:2:1171:A:H2'	1:2:1172:G:C8	2.39	0.58
1:2:649:U:H2'	1:2:650:U:C6	2.40	0.57
1:2:828:U:H2'	1:2:829:A:C8	2.39	0.57
1:2:1160:A:H2'	1:2:1161:C:C6	2.39	0.57
1:2:12:U:H2'	1:2:13:C:C6	2.40	0.57
1:2:1171:A:H2'	1:2:1172:G:H8	1.70	0.57
1:2:183:U:H2'	1:2:184:C:C6	2.40	0.57
11:T:39:GLU:HB3	11:T:46:LYS:HA	1.87	0.57
1:2:650:U:H2'	1:2:651:G:C8	2.39	0.56
1:2:1749:A:C3'	1:2:1750:A:H4'	2.35	0.56
1:2:1375:A:H2'	1:2:1376:C:C6	2.41	0.56
1:2:258:C:H2'	1:2:259:U:H5'	1.88	0.56
24:p:120:LEU:HD23	24:p:128:VAL:HG11	1.87	0.56
20:d:29:HIS:O	20:d:29:HIS:ND1	2.39	0.55
26:t:294:HIS:HE1	26:t:296:PRO:HA	1.71	0.55
8:L:27:GLN:OE1	8:L:65:ARG:NH2	2.34	0.55
1:2:1253:U:H2'	1:2:1254:U:C6	2.41	0.55
25:r:86:LYS:HE2	25:r:87:THR:HG23	1.88	0.55
1:2:1534:G:OP2	7:K:74:SER:OG	2.21	0.55
7:K:95:HIS:O	7:K:97:LYS:N	2.40	0.54
1:2:330:G:OP2	13:V:172:ARG:NH1	2.40	0.54
25:r:185:HIS:O	25:r:185:HIS:ND1	2.41	0.54
1:2:906:A:H2'	1:2:907:A:C8	2.43	0.54
9:Q:176:VAL:HG12	9:Q:177:GLN:H	1.72	0.54
1:2:127:G:N2	1:2:178:U:O2'	2.40	0.54
18:b:12:ASN:HD22	18:b:12:ASN:C	2.16	0.54
1:2:181:A:H2'	1:2:182:A:C8	2.43	0.54
1:2:198:A:H2'	1:2:199:G:O4'	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1474:G:H2'	1:2:1475:A:C8	2.43	0.53
9:Q:41:ARG:HH21	9:Q:232:HIS:CD2	2.26	0.53
19:c:65:ASN:ND2	19:c:116:ASP:OD2	2.31	0.53
25:r:88:MET:HA	25:r:91:ARG:HG2	1.89	0.53
1:2:184:C:H2'	1:2:185:U:C6	2.42	0.53
1:2:184:C:H2'	1:2:185:U:H6	1.74	0.53
1:2:555:A:H8	1:2:590:C:O2'	1.91	0.53
1:2:850:A:H2'	1:2:851:U:C6	2.42	0.53
1:2:1564:U:H2'	1:2:1565:C:H6	1.73	0.53
26:t:267:VAL:O	26:t:275:ASP:N	2.33	0.53
1:2:821:U:H3	1:2:852:C:N4	2.07	0.53
1:2:1348:A:H2'	1:2:1349:G:C8	2.44	0.53
1:2:1349:G:H2'	1:2:1350:U:C6	2.44	0.53
1:2:1474:G:H2'	1:2:1475:A:H8	1.73	0.53
1:2:72:A:N7	11:T:167:LYS:NZ	2.51	0.53
1:2:1374:C:H2'	1:2:1375:A:C8	2.43	0.53
1:2:183:U:H2'	1:2:184:C:H6	1.74	0.52
1:2:1452:U:H2'	1:2:1453:G:H8	1.75	0.52
26:t:484:ASP:OD1	26:t:644:ARG:NH1	2.42	0.52
1:2:520:A:H2'	1:2:521:A:C8	2.45	0.52
1:2:885:G:H2'	1:2:886:U:C6	2.44	0.52
1:2:1656:U:C2	1:2:1658:G:C8	2.98	0.52
1:2:1590:G:H2'	1:2:1591:C:H6	1.75	0.52
26:t:638:ASN:ND2	26:t:694:SER:OG	2.37	0.52
1:2:250:C:H5'	1:2:250:C:C6	2.44	0.52
26:t:237:ASN:C	26:t:237:ASN:HD22	2.16	0.52
8:L:44:VAL:HG22	8:L:54:LEU:HD21	1.92	0.52
26:t:514:VAL:O	26:t:776:ARG:NH2	2.31	0.51
26:t:595:ASN:ND2	26:t:690:ASN:OD1	2.36	0.51
1:2:1588:G:O6	1:2:1608:U:O4	2.29	0.51
1:2:555:A:H8	1:2:590:C:HO2'	1.53	0.51
1:2:1173:C:H2'	1:2:1174:C:C6	2.46	0.51
1:2:830:U:H2'	1:2:831:U:O4'	2.10	0.51
1:2:1748:G:H2'	1:2:1749:A:N7	2.26	0.51
1:2:1508:U:H2'	1:2:1509:C:C6	2.46	0.50
26:t:294:HIS:HB2	26:t:584:LEU:HD21	1.94	0.50
1:2:139:C:OP2	11:T:187:LYS:NZ	2.44	0.50
1:2:895:G:H2'	1:2:896:U:C6	2.45	0.50
12:U:15:GLU:OE1	12:U:19:GLN:NE2	2.36	0.50
1:2:1220:C:H2'	1:2:1221:A:C8	2.46	0.50
1:2:1564:U:H2'	1:2:1565:C:C6	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1350:U:H2'	1:2:1351:G:C8	2.46	0.50
25:r:125:HIS:NE2	25:r:162:GLU:OE1	2.24	0.50
1:2:1688:U:O4	1:2:1713:G:N2	2.40	0.50
26:t:294:HIS:CE1	26:t:296:PRO:HA	2.46	0.50
1:2:182:A:H2'	1:2:183:U:C6	2.45	0.50
1:2:1055:U:H2'	1:2:1056:U:C6	2.47	0.50
1:2:1579:U:H2'	1:2:1580:C:C6	2.46	0.49
1:2:1690:G:C2	1:2:1712:A:C6	3.00	0.49
19:c:90:ASP:CG	19:c:91:GLY:H	2.18	0.49
1:2:649:U:H2'	1:2:650:U:H6	1.75	0.49
1:2:1565:C:OP1	5:H:41:ARG:HD3	2.13	0.49
1:2:650:U:H2'	1:2:651:G:H8	1.77	0.49
2:B:62:VAL:HG13	2:B:89:ILE:HG13	1.94	0.49
1:2:1452:U:H2'	1:2:1453:G:C8	2.48	0.49
7:K:59:TYR:HE1	7:K:100:ILE:HG23	1.78	0.49
26:t:600:ARG:NH1	26:t:604:TYR:O	2.44	0.49
1:2:846:G:N3	1:2:846:G:H5'	2.27	0.49
1:2:1477:G:H2'	1:2:1478:G:H8	1.78	0.49
1:2:1503:A:H2'	1:2:1504:G:C8	2.48	0.49
19:c:73:ARG:HG3	19:c:73:ARG:HH11	1.78	0.49
1:2:1656:U:H2'	1:2:1657:U:H5'	1.95	0.48
1:2:853:G:OP2	1:2:853:G:C8	2.66	0.48
1:2:1241:G:H1'	3:E:79:HIS:NE2	2.28	0.48
1:2:1533:C:OP2	7:K:77:ARG:NH2	2.39	0.48
24:p:204:ARG:NH1	24:p:250:LEU:O	2.39	0.48
1:2:284:G:N2	11:T:188:ARG:HD2	2.29	0.48
1:2:514:G:HO2'	1:2:515:A:H8	1.62	0.48
1:2:1590:G:H2'	1:2:1591:C:C6	2.48	0.48
2:B:225:ARG:O	8:L:61:ARG:NH2	2.47	0.48
1:2:1238:A:H2'	1:2:1239:U:C5	2.49	0.48
1:2:1461:C:H2'	1:2:1462:G:H8	1.78	0.48
1:2:1647:U:H2'	1:2:1648:A:C8	2.49	0.48
2:B:161:ASP:O	8:L:45:LYS:N	2.45	0.48
1:2:992:A:H2'	1:2:993:A:H5'	1.96	0.48
1:2:821:U:H2'	1:2:822:U:C6	2.49	0.47
1:2:1691:A:N6	1:2:1710:U:OP2	2.44	0.47
1:2:1749:A:H3'	1:2:1750:A:H4'	1.95	0.47
14:W:53:ARG:NH2	14:W:99:LEU:O	2.47	0.47
1:2:868:G:H1	1:2:960:U:H3	1.60	0.47
2:B:25:LEU:HD22	4:F:27:GLY:HA3	1.95	0.47
11:T:78:THR:O	11:T:81:VAL:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:U:75:THR:O	12:U:76:LYS:HB2	2.14	0.47
1:2:1713:G:H2'	1:2:1714:A:C8	2.49	0.47
1:2:1489:U:H2'	1:2:1514:U:O4	2.14	0.47
1:2:1689:A:C8	1:2:1690:G:N2	2.82	0.47
1:2:1241:G:H5'	3:E:77:ARG:HG3	1.95	0.47
1:2:1357:A:H2'	1:2:1358:G:C8	2.50	0.47
1:2:12:U:H2'	1:2:13:C:H6	1.80	0.47
1:2:848:C:H2'	1:2:849:C:H6	1.79	0.47
1:2:929:A:C8	17:Z:123:SER:O	2.68	0.47
1:2:1071:U:H2'	1:2:1072:C:C6	2.50	0.47
1:2:1613:U:OP1	2:B:169:ASN:ND2	2.48	0.47
1:2:1615:C:O2'	1:2:1616:G:OP2	2.26	0.47
16:Y:3:ARG:NH2	16:Y:10:GLY:O	2.48	0.47
1:2:250:C:H5'	1:2:250:C:H6	1.80	0.47
1:2:281:G:O2'	1:2:282:C:OP2	2.26	0.47
1:2:852:C:H2'	1:2:853:G:O4'	2.15	0.47
1:2:555:A:C8	1:2:590:C:O2'	2.66	0.47
1:2:560:U:H2'	1:2:561:G:H8	1.80	0.47
25:r:223:SER:HB3	25:r:321:LEU:HD12	1.96	0.47
1:2:852:C:O2'	1:2:853:G:H5'	2.15	0.46
1:2:1206:U:H3'	1:2:1207:C:H2'	1.96	0.46
9:Q:78:ASP:OD1	9:Q:79:HIS:ND1	2.46	0.46
16:Y:87:ASP:OD1	16:Y:88:LEU:N	2.48	0.46
1:2:108:A:H2'	1:2:109:G:C8	2.50	0.46
19:c:68:ILE:O	19:c:70:LYS:NZ	2.41	0.46
1:2:1374:C:H2'	1:2:1375:A:H8	1.80	0.46
12:U:15:GLU:HA	12:U:18:LEU:HD12	1.97	0.46
14:W:163:PRO:C	14:W:165:GLY:H	2.23	0.46
1:2:1202:A:H1'	1:2:1207:C:H42	1.81	0.46
1:2:1357:A:H2'	1:2:1358:G:H8	1.80	0.46
2:B:183:ALA:O	2:B:186:ASN:HB3	2.15	0.46
26:t:601:TRP:CD1	26:t:602:GLU:H	2.33	0.46
14:W:117:GLY:O	14:W:119:ALA:N	2.49	0.46
1:2:823:G:O6	1:2:850:A:N6	2.49	0.46
1:2:1161:C:H2'	1:2:1162:C:H6	1.80	0.46
1:2:871:G:H2'	1:2:872:G:C8	2.51	0.46
1:2:1523:G:N3	1:2:1523:G:H5'	2.31	0.46
1:2:1659:A:O2'	1:2:1660:A:O5'	2.33	0.46
12:U:28:GLU:HG3	12:U:35:LYS:HA	1.97	0.46
1:2:647:G:H2'	1:2:648:G:C8	2.51	0.46
1:2:826:U:H2'	1:2:827:C:C6	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1350:U:H2'	1:2:1351:G:H8	1.81	0.46
9:Q:179:SER:HB3	9:Q:183:GLN:NE2	2.31	0.46
26:t:84:LYS:NZ	26:t:242:LEU:O	2.33	0.46
1:2:853:G:OP2	1:2:853:G:H8	1.99	0.46
19:c:59:ILE:O	19:c:69:ARG:N	2.43	0.46
1:2:684:A:H2'	1:2:685:A:C8	2.50	0.45
1:2:67:A:H2'	1:2:69:G:H5''	1.98	0.45
1:2:116:U:H2'	1:2:117:U:C6	2.51	0.45
1:2:892:A:H2'	1:2:893:U:H6	1.81	0.45
1:2:1082:C:H3'	1:2:1083:G:H5''	1.98	0.45
1:2:939:A:H2'	1:2:940:A:C8	2.51	0.45
7:K:61:SER:H	7:K:64:VAL:HB	1.82	0.45
1:2:1239:U:C2	1:2:1243:G:N7	2.85	0.45
1:2:1651:A:H2'	1:2:1652:C:C6	2.51	0.45
6:I:35:ASP:C	6:I:35:ASP:OD1	2.60	0.45
12:U:134:GLU:OE1	16:Y:21:ASN:ND2	2.40	0.45
1:2:1652:C:H2'	1:2:1653:C:C6	2.51	0.45
1:2:1711:C:H4'	1:2:1712:A:OP2	2.16	0.45
1:2:1458:G:N3	1:2:1458:G:H2'	2.31	0.45
1:2:1066:C:H4'	9:Q:149:GLN:HG2	1.98	0.45
3:E:83:MET:HE2	3:E:89:MET:HE1	1.98	0.45
1:2:1358:G:O2'	6:I:133:ASP:OD2	2.28	0.45
1:2:1714:A:H2'	1:2:1715:G:C8	2.51	0.45
1:2:198:A:O5'	1:2:198:A:H8	2.00	0.45
9:Q:109:LYS:HE3	9:Q:113:MET:HE3	1.99	0.45
11:T:1:MET:HE1	11:T:106:LEU:HB2	1.98	0.44
1:2:856:A:N6	12:U:96:ARG:HB3	2.32	0.44
1:2:1539:G:H5'	1:2:1539:G:C8	2.52	0.44
1:2:848:C:H2'	1:2:849:C:C6	2.52	0.44
1:2:1741:U:H2'	1:2:1742:U:C6	2.52	0.44
1:2:892:A:H2'	1:2:893:U:C6	2.53	0.44
1:2:895:G:H21	17:Z:38:THR:HG21	1.82	0.44
1:2:1248:C:C2	1:2:1249:U:C5	3.05	0.44
1:2:1713:G:H2'	1:2:1714:A:H8	1.81	0.44
1:2:1494:C:H2'	1:2:1495:C:C6	2.52	0.44
1:2:1749:A:H2'	1:2:1750:A:H4'	2.00	0.44
1:2:953:G:H2'	1:2:954:G:C8	2.53	0.44
1:2:1173:C:H2'	1:2:1174:C:H6	1.81	0.44
1:2:1749:A:O5'	1:2:1749:A:H8	2.00	0.44
12:U:91:ILE:HG12	12:U:129:LEU:HD23	2.00	0.44
13:V:151:LYS:HE2	13:V:151:LYS:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:821:U:O2	1:2:852:C:N3	2.51	0.44
1:2:1186:U:O4	1:2:1208:A:N6	2.50	0.44
1:2:1451:C:H2'	1:2:1452:U:C6	2.52	0.44
1:2:844:A:H2'	1:2:845:G:C8	2.53	0.44
1:2:1606:C:H2'	1:2:1607:G:C8	2.52	0.44
1:2:1003:A:N6	24:p:242:MET:HG2	2.33	0.43
10:S:139:VAL:HG13	10:S:150:PRO:HG3	2.00	0.43
26:t:188:TYR:HE1	26:t:223:GLU:HG2	1.83	0.43
1:2:591:A:H2'	1:2:592:A:C8	2.53	0.43
1:2:1632:C:OP2	24:p:264:ARG:NH1	2.51	0.43
1:2:1750:A:N3	1:2:1750:A:H2'	2.34	0.43
9:Q:28:GLU:N	9:Q:48:VAL:O	2.48	0.43
16:Y:103:GLU:O	16:Y:106:ARG:NH1	2.49	0.43
1:2:482:U:H2'	1:2:483:A:H8	1.83	0.43
1:2:1477:G:H2'	1:2:1478:G:C8	2.53	0.43
1:2:1489:U:C5	1:2:1513:G:C6	3.06	0.43
1:2:1749:A:C8	1:2:1750:A:H1'	2.53	0.43
1:2:1558:U:H5'	5:H:126:ARG:NH2	2.33	0.43
20:d:124:ARG:O	20:d:125:LEU:HB3	2.18	0.43
1:2:894:U:H2'	1:2:895:G:C8	2.54	0.43
1:2:1568:C:O2'	1:2:1569:A:OP2	2.18	0.43
1:2:1652:C:H2'	1:2:1653:C:C1'	2.49	0.43
10:S:100:ARG:HH21	10:S:118:GLU:HG2	1.81	0.43
21:e:258:VAL:O	21:e:262:ALA:N	2.37	0.43
1:2:1654:G:H1'	1:2:1746:A:H61	1.84	0.43
19:c:27:ASN:C	19:c:27:ASN:HD22	2.27	0.43
1:2:198:A:O5'	1:2:198:A:C8	2.71	0.43
1:2:1655:A:H2'	1:2:1656:U:C6	2.54	0.43
2:B:149:VAL:HG12	2:B:156:ARG:H	1.83	0.43
26:t:259:VAL:HG12	26:t:282:VAL:HA	2.01	0.43
1:2:126:A:OP1	11:T:201:GLN:HG3	2.18	0.43
1:2:845:G:H2'	1:2:846:G:C5'	2.48	0.43
1:2:1259:U:H2'	1:2:1260:U:C6	2.54	0.43
10:S:94:ALA:C	10:S:96:ASN:H	2.27	0.42
1:2:843:U:O4	1:2:844:A:N6	2.51	0.42
1:2:849:C:H2'	1:2:850:A:H8	1.83	0.42
1:2:1077:C:H2'	1:2:1078:C:H6	1.83	0.42
1:2:1749:A:N7	1:2:1750:A:C8	2.87	0.42
11:T:78:THR:HG22	11:T:79:LYS:N	2.34	0.42
1:2:221:A:C8	1:2:833:U:H1'	2.53	0.42
1:2:260:U:H3'	1:2:261:U:H5'	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:U:109:VAL:HG12	12:U:110:GLN:N	2.34	0.42
1:2:545:A:H2'	23:g:31:LYS:CD	2.48	0.42
5:H:83:ALA:HA	5:H:86:LEU:HD13	2.02	0.42
1:2:540:G:O2'	1:2:541:A:OP1	2.37	0.42
1:2:1568:C:OP1	5:H:36:LYS:NZ	2.51	0.42
1:2:647:G:H21	1:2:687:G:H1	1.67	0.42
1:2:819:G:C6	1:2:853:G:C5	3.07	0.42
1:2:842:C:H2'	1:2:843:U:O4'	2.20	0.42
1:2:1622:G:H2'	1:2:1623:C:H6	1.85	0.42
18:b:6:VAL:HG12	18:b:34:ILE:HD11	2.02	0.42
26:t:493:SER:OG	26:t:496:GLU:OE1	2.24	0.42
1:2:1677:C:OP1	13:V:42:ARG:NH1	2.50	0.42
11:T:185:GLN:HA	11:T:188:ARG:NH1	2.35	0.42
1:2:830:U:C2'	1:2:831:U:H5'	2.49	0.41
4:F:69:VAL:HG21	4:F:77:GLN:HB2	2.02	0.41
1:2:1260:U:H2'	1:2:1261:G:H8	1.85	0.41
1:2:1492:A:O2'	1:2:1493:A:P	2.79	0.41
25:r:204:LYS:H	25:r:204:LYS:HD3	1.84	0.41
26:t:482:PHE:HB3	26:t:485:GLU:CD	2.45	0.41
1:2:819:G:C4	1:2:853:G:C2	3.08	0.41
1:2:897:C:O2'	1:2:914:G:N2	2.53	0.41
1:2:959:U:O2	1:2:959:U:H2'	2.19	0.41
26:t:44:LYS:HB2	26:t:45:PRO:HD3	2.02	0.41
26:t:522:PRO:HD2	26:t:525:TRP:HE1	1.85	0.41
26:t:708:PHE:CD2	26:t:709:ARG:HD3	2.56	0.41
1:2:845:G:H2'	1:2:846:G:H5''	2.03	0.41
5:H:96:LYS:HB2	5:H:98:TYR:CE1	2.55	0.41
24:p:169:ASP:OD1	24:p:169:ASP:N	2.49	0.41
1:2:1157:A:H2'	1:2:1160:A:N7	2.36	0.41
1:2:1568:C:HO2'	1:2:1569:A:P	2.39	0.41
1:2:1603:U:H2'	1:2:1604:U:C6	2.56	0.41
1:2:821:U:C2	1:2:852:C:N3	2.88	0.41
5:H:94:ASP:C	5:H:96:LYS:H	2.28	0.41
9:Q:104:ASP:OD1	9:Q:105:PHE:N	2.53	0.41
11:T:116:LYS:HE3	11:T:125:THR:HG21	2.01	0.41
22:f:37:CYS:SG	22:f:40:CYS:N	2.94	0.41
1:2:182:A:H2'	1:2:183:U:H6	1.84	0.41
1:2:819:G:H1'	1:2:820:U:OP2	2.21	0.41
1:2:1202:A:H1'	1:2:1207:C:N4	2.36	0.41
1:2:219:A:N6	1:2:843:U:C2	2.88	0.41
1:2:639:U:OP1	12:U:112:ARG:NH2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:851:U:H2'	1:2:852:C:C5	2.56	0.41
1:2:1511:U:H2'	1:2:1512:G:C8	2.56	0.41
1:2:1690:G:C2	1:2:1712:A:N6	2.88	0.41
2:B:216:GLU:OE1	2:B:219:ARG:NH1	2.47	0.41
5:H:88:ARG:HB2	5:H:98:TYR:O	2.21	0.41
14:W:65:LYS:HA	14:W:70:LEU:HD11	2.02	0.41
1:2:197:A:O2'	1:2:198:A:O4'	2.21	0.41
1:2:1461:C:H2'	1:2:1462:G:C8	2.55	0.41
1:2:1691:A:C2	1:2:1710:U:H6	2.38	0.41
3:E:19:GLY:N	5:H:93:THR:O	2.52	0.41
26:t:265:ASP:O	26:t:276:LEU:HD12	2.20	0.41
1:2:1483:A:N3	1:2:1607:G:O2'	2.44	0.41
19:c:41:SER:HA	19:c:42:PRO:HD3	1.92	0.41
25:r:322:ASP:N	25:r:322:ASP:OD1	2.54	0.41
1:2:514:G:O2'	1:2:515:A:H8	2.03	0.40
1:2:562:G:H2'	1:2:563:U:C6	2.56	0.40
11:T:20:ASP:HB2	11:T:23:ARG:HE	1.86	0.40
1:2:560:U:H2'	1:2:561:G:C8	2.56	0.40
1:2:755:A:H8	1:2:755:A:OP2	2.05	0.40
1:2:181:A:H2'	1:2:182:A:H8	1.87	0.40
1:2:1469:A:H2'	1:2:1470:C:C6	2.57	0.40
14:W:70:LEU:HD23	14:W:70:LEU:HA	1.91	0.40
1:2:1003:A:H62	24:p:242:MET:HG2	1.86	0.40
5:H:81:ILE:HA	5:H:82:PRO:HD3	1.93	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	197/225 (88%)	189 (96%)	7 (4%)	1 (0%)	24 54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	E	110/142 (78%)	102 (93%)	8 (7%)	0	100	100
4	F	123/143 (86%)	115 (94%)	7 (6%)	1 (1%)	16	44
5	H	120/146 (82%)	110 (92%)	10 (8%)	0	100	100
6	I	139/144 (96%)	136 (98%)	2 (1%)	1 (1%)	18	48
7	K	67/108 (62%)	64 (96%)	2 (3%)	1 (2%)	8	28
8	L	58/67 (87%)	55 (95%)	3 (5%)	0	100	100
9	Q	212/255 (83%)	208 (98%)	4 (2%)	0	100	100
10	S	258/261 (99%)	250 (97%)	8 (3%)	0	100	100
11	T	224/236 (95%)	219 (98%)	4 (2%)	1 (0%)	30	59
12	U	182/190 (96%)	172 (94%)	9 (5%)	1 (0%)	24	54
13	V	181/200 (90%)	175 (97%)	6 (3%)	0	100	100
14	W	183/197 (93%)	178 (97%)	4 (2%)	1 (0%)	24	54
15	X	139/156 (89%)	135 (97%)	4 (3%)	0	100	100
16	Y	148/151 (98%)	143 (97%)	5 (3%)	0	100	100
17	Z	125/137 (91%)	119 (95%)	4 (3%)	2 (2%)	7	27
18	b	127/130 (98%)	123 (97%)	4 (3%)	0	100	100
19	c	142/145 (98%)	134 (94%)	8 (6%)	0	100	100
20	d	128/135 (95%)	124 (97%)	4 (3%)	0	100	100
21	e	167/483 (35%)	162 (97%)	4 (2%)	1 (1%)	21	51
22	f	79/82 (96%)	77 (98%)	1 (1%)	1 (1%)	9	32
23	g	26/63 (41%)	24 (92%)	2 (8%)	0	100	100
24	p	178/274 (65%)	174 (98%)	4 (2%)	0	100	100
25	r	174/425 (41%)	166 (95%)	8 (5%)	0	100	100
26	t	607/788 (77%)	587 (97%)	19 (3%)	1 (0%)	43	72
All	All	4094/5283 (78%)	3941 (96%)	141 (3%)	12 (0%)	37	65

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	K	96	SER
17	Z	42	VAL
12	U	98	ILE
14	W	118	LEU
26	t	96	ASP

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Mol	Chain	Res	Type
17	Z	13	VAL
4	F	97	VAL
22	f	60	SER
21	e	282	ILE
2	B	23	VAL
11	T	70	PRO
6	I	45	MET

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	173/191 (91%)	171 (99%)	2 (1%)	63	86
3	E	95/118 (80%)	95 (100%)	0	100	100
4	F	105/119 (88%)	103 (98%)	2 (2%)	50	79
5	H	109/129 (84%)	108 (99%)	1 (1%)	70	90
6	I	113/116 (97%)	110 (97%)	3 (3%)	39	73
7	K	60/89 (67%)	60 (100%)	0	100	100
8	L	53/60 (88%)	53 (100%)	0	100	100
9	Q	191/224 (85%)	191 (100%)	0	100	100
10	S	221/222 (100%)	221 (100%)	0	100	100
11	T	191/201 (95%)	189 (99%)	2 (1%)	68	89
12	U	165/170 (97%)	163 (99%)	2 (1%)	63	86
13	V	147/161 (91%)	146 (99%)	1 (1%)	76	92
14	W	158/166 (95%)	158 (100%)	0	100	100
15	X	126/137 (92%)	124 (98%)	2 (2%)	55	83
16	Y	127/128 (99%)	127 (100%)	0	100	100
17	Z	91/105 (87%)	90 (99%)	1 (1%)	65	88
18	b	110/111 (99%)	109 (99%)	1 (1%)	70	90
19	c	119/120 (99%)	118 (99%)	1 (1%)	73	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
20	d	109/113 (96%)	108 (99%)	1 (1%)	70	90
21	e	17/424 (4%)	17 (100%)	0	100	100
22	f	70/71 (99%)	69 (99%)	1 (1%)	59	85
23	g	27/54 (50%)	27 (100%)	0	100	100
24	p	157/238 (66%)	154 (98%)	3 (2%)	50	79
25	r	171/384 (44%)	169 (99%)	2 (1%)	63	86
26	t	535/703 (76%)	531 (99%)	4 (1%)	76	92
All	All	3440/4554 (76%)	3411 (99%)	29 (1%)	70	90

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

Mol	Chain	Res	Type
2	B	76	ARG
2	B	122	ASN
4	F	70	THR
4	F	115	THR
5	H	17	LEU
6	I	33	TYR
6	I	35	ASP
6	I	95	ASP
11	T	23	ARG
11	T	168	THR
12	U	14	THR
12	U	73	VAL
13	V	67	TRP
15	X	98	ASN
15	X	109	VAL
17	Z	42	VAL
18	b	12	ASN
19	c	60	GLU
20	d	121	THR
22	f	43	ILE
24	p	116	ILE
24	p	145	PHE
24	p	245	GLU
25	r	152	MET
25	r	182	ASN
26	t	154	VAL
26	t	191	VAL
26	t	338	GLU

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Mol	Chain	Res	Type
26	t	347	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	169	ASN
2	B	200	ASN
4	F	8	GLN
4	F	21	HIS
4	F	74	HIS
5	H	13	HIS
5	H	75	ASN
5	H	78	HIS
6	I	23	GLN
6	I	77	ASN
6	I	93	HIS
7	K	44	GLN
7	K	95	HIS
9	Q	95	ASN
9	Q	157	GLN
9	Q	177	GLN
10	S	153	ASN
10	S	201	HIS
10	S	258	GLN
11	T	10	ASN
12	U	29	ASN
12	U	161	GLN
13	V	116	HIS
14	W	48	GLN
14	W	112	GLN
16	Y	58	HIS
17	Z	80	HIS
18	b	12	ASN
18	b	15	ASN
18	b	24	GLN
18	b	56	HIS
19	c	27	ASN
20	d	107	GLN
20	d	110	GLN
22	f	9	HIS
26	t	14	HIS
26	t	134	ASN

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Mol	Chain	Res	Type
26	t	205	GLN
26	t	237	ASN
26	t	294	HIS
26	t	340	ASN
26	t	572	GLN
26	t	589	HIS
26	t	640	HIS
26	t	711	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	2	1397/1800 (77%)	177 (12%)	51 (3%)

All (177) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	2	3	U
1	2	4	C
1	2	26	A
1	2	34	G
1	2	42	G
1	2	47	A
1	2	48	G
1	2	68	A
1	2	69	G
1	2	72	A
1	2	77	U
1	2	100	A
1	2	104	A
1	2	114	C
1	2	116	U
1	2	127	G
1	2	129	U
1	2	139	C
1	2	140	A
1	2	144	U
1	2	146	U
1	2	159	U
1	2	161	U
1	2	204	G

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Mol	Chain	Res	Type
1	2	220	A
1	2	249	U
1	2	250	C
1	2	251	A
1	2	259	U
1	2	260	U
1	2	261	U
1	2	265	A
1	2	276	C
1	2	282	C
1	2	314	C
1	2	316	A
1	2	321	C
1	2	322	G
1	2	323	A
1	2	337	G
1	2	338	C
1	2	359	A
1	2	360	A
1	2	361	C
1	2	400	A
1	2	401	A
1	2	402	C
1	2	404	G
1	2	416	A
1	2	424	C
1	2	426	G
1	2	439	U
1	2	444	C
1	2	453	U
1	2	454	U
1	2	468	A
1	2	469	C
1	2	487	G
1	2	510	G
1	2	511	A
1	2	514	G
1	2	515	A
1	2	540	G
1	2	541	A
1	2	542	A
1	2	557	G

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Mol	Chain	Res	Type
1	2	558	U
1	2	559	C
1	2	560	U
1	2	577	G
1	2	594	A
1	2	595	G
1	2	606	A
1	2	608	U
1	2	620	A
1	2	622	A
1	2	623	A
1	2	639	U
1	2	694	U
1	2	765	G
1	2	766	U
1	2	774	A
1	2	789	A
1	2	811	A
1	2	812	A
1	2	813	U
1	2	814	A
1	2	815	G
1	2	819	G
1	2	820	U
1	2	821	U
1	2	831	U
1	2	833	U
1	2	840	U
1	2	846	G
1	2	853	G
1	2	856	A
1	2	863	A
1	2	886	U
1	2	898	A
1	2	914	G
1	2	933	A
1	2	935	U
1	2	942	G
1	2	960	U
1	2	966	A
1	2	992	A
1	2	993	A

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Mol	Chain	Res	Type
1	2	1000	C
1	2	1005	A
1	2	1025	A
1	2	1028	C
1	2	1031	U
1	2	1040	G
1	2	1043	A
1	2	1052	U
1	2	1053	G
1	2	1058	U
1	2	1083	G
1	2	1092	A
1	2	1096	C
1	2	1097	U
1	2	1138	A
1	2	1158	C
1	2	1159	C
1	2	1160	A
1	2	1185	U
1	2	1207	C
1	2	1214	U
1	2	1242	A
1	2	1243	G
1	2	1244	A
1	2	1245	G
1	2	1471	A
1	2	1491	U
1	2	1492	A
1	2	1493	A
1	2	1515	A
1	2	1516	A
1	2	1523	G
1	2	1524	A
1	2	1535	U
1	2	1536	G
1	2	1537	C
1	2	1538	U
1	2	1557	U
1	2	1560	U
1	2	1568	C
1	2	1569	A
1	2	1584	G

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Mol	Chain	Res	Type
1	2	1595	U
1	2	1596	C
1	2	1616	G
1	2	1630	U
1	2	1631	A
1	2	1632	C
1	2	1650	U
1	2	1653	C
1	2	1657	U
1	2	1659	A
1	2	1660	A
1	2	1688	U
1	2	1689	A
1	2	1690	G
1	2	1691	A
1	2	1711	C
1	2	1712	A
1	2	1715	G
1	2	1745	G
1	2	1747	G
1	2	1750	A
1	2	1751	C
1	2	1770	U
1	2	1780	G
1	2	1781	A
1	2	1792	G
1	2	1795	U

All (51) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	2	3	U
1	2	9	U
1	2	47	A
1	2	138	A
1	2	249	U
1	2	250	C
1	2	315	A
1	2	321	C
1	2	322	G
1	2	400	A
1	2	401	A

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Mol	Chain	Res	Type
1	2	453	U
1	2	468	A
1	2	510	G
1	2	514	G
1	2	557	G
1	2	558	U
1	2	606	A
1	2	622	A
1	2	765	G
1	2	811	A
1	2	814	A
1	2	819	G
1	2	855	A
1	2	999	U
1	2	1052	U
1	2	1057	U
1	2	1092	A
1	2	1158	C
1	2	1241	G
1	2	1243	G
1	2	1364	G
1	2	1491	U
1	2	1492	A
1	2	1535	U
1	2	1537	C
1	2	1556	A
1	2	1568	C
1	2	1595	U
1	2	1615	C
1	2	1630	U
1	2	1631	A
1	2	1649	G
1	2	1657	U
1	2	1659	A
1	2	1688	U
1	2	1710	U
1	2	1711	C
1	2	1744	A
1	2	1750	A
1	2	1780	G

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

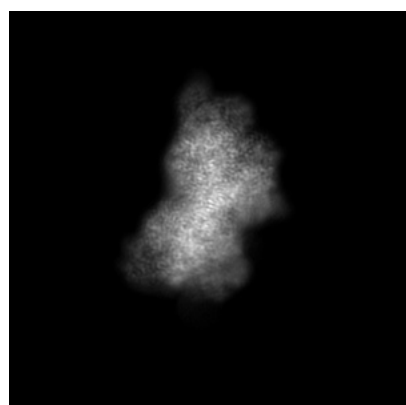
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-16347. These allow visual inspection of the internal detail of the map and identification of artifacts.

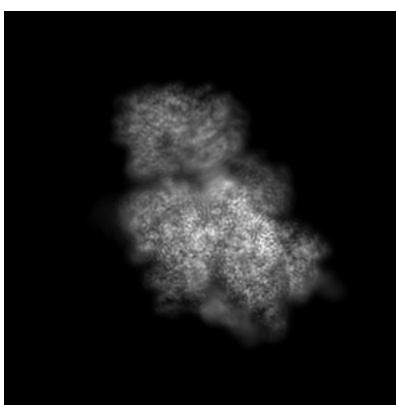
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

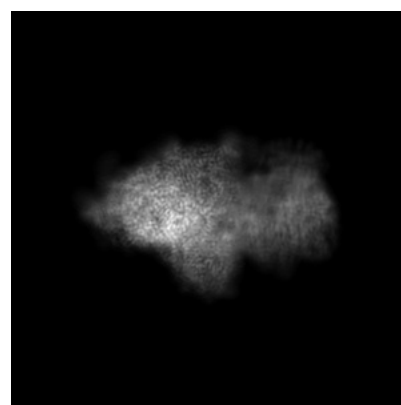
6.1.1 Primary map



X



Y

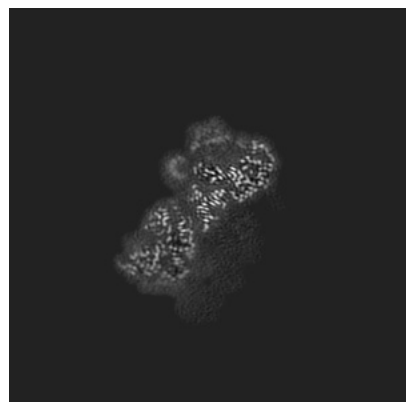


Z

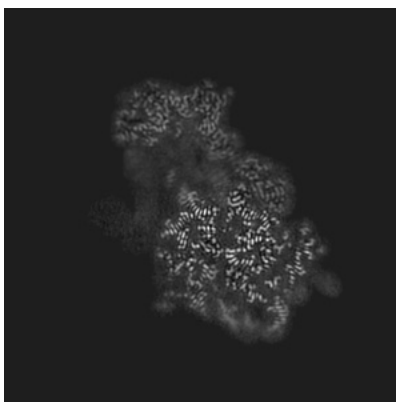
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

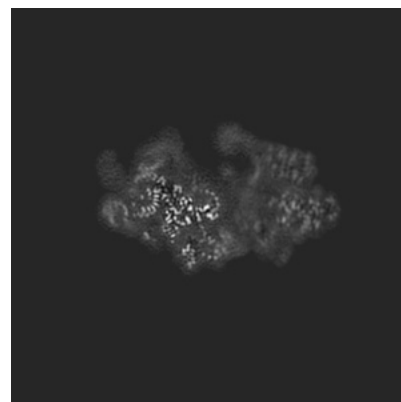
6.2.1 Primary map



X Index: 200



Y Index: 200

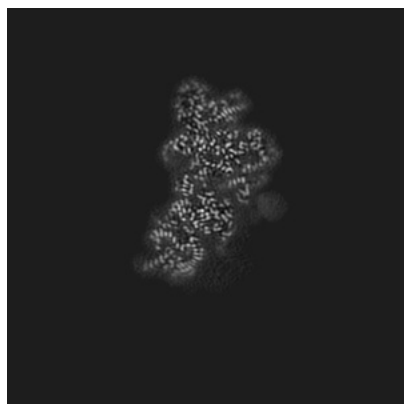


Z Index: 200

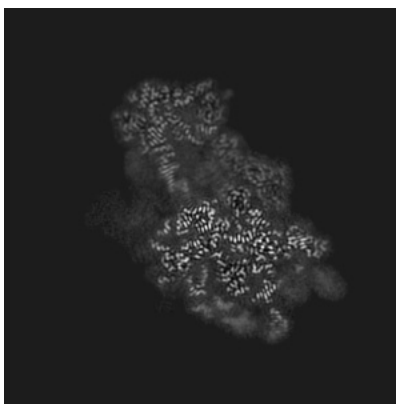
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

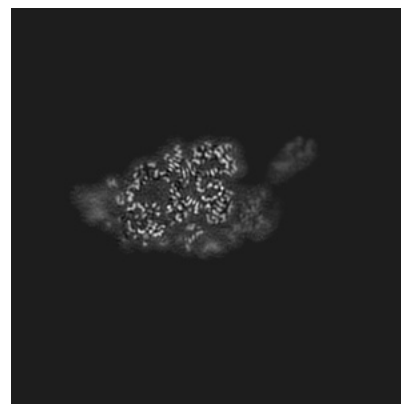
6.3.1 Primary map



X Index: 161



Y Index: 193

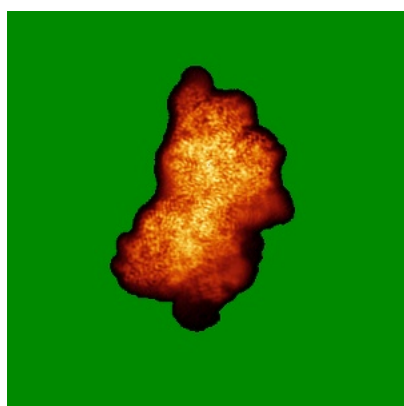


Z Index: 240

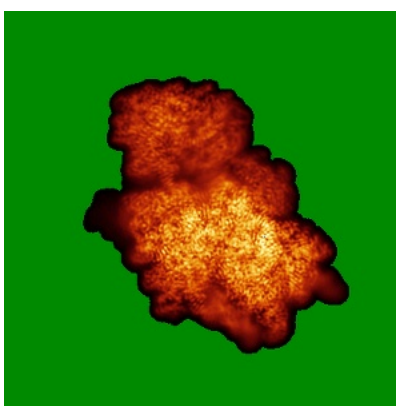
The images above show the largest variance slices of the map in three orthogonal directions.

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

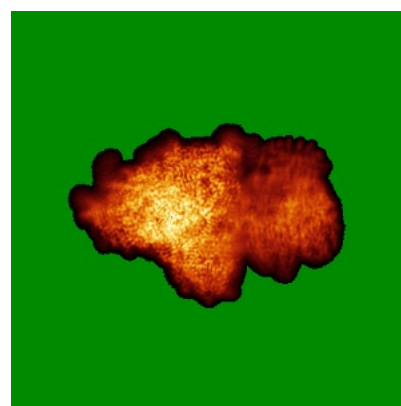
6.4.1 Primary map



X



Y

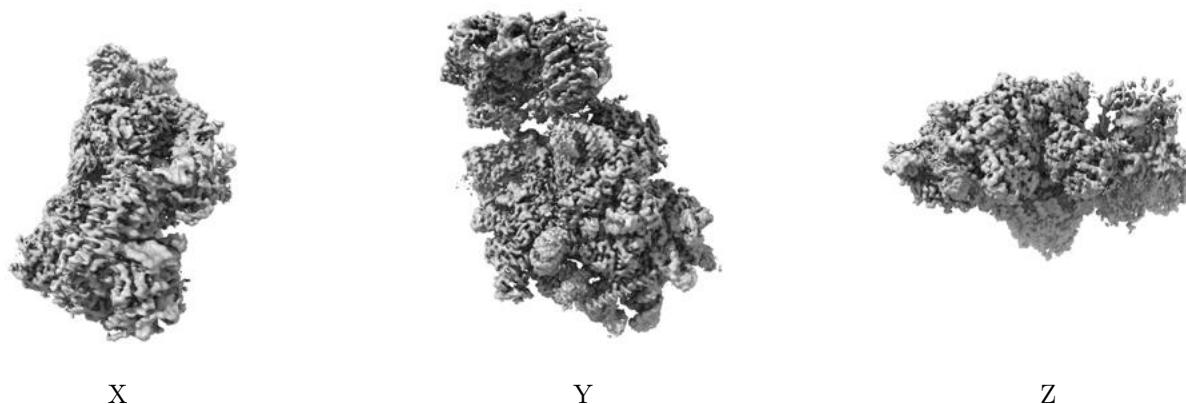


Z

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

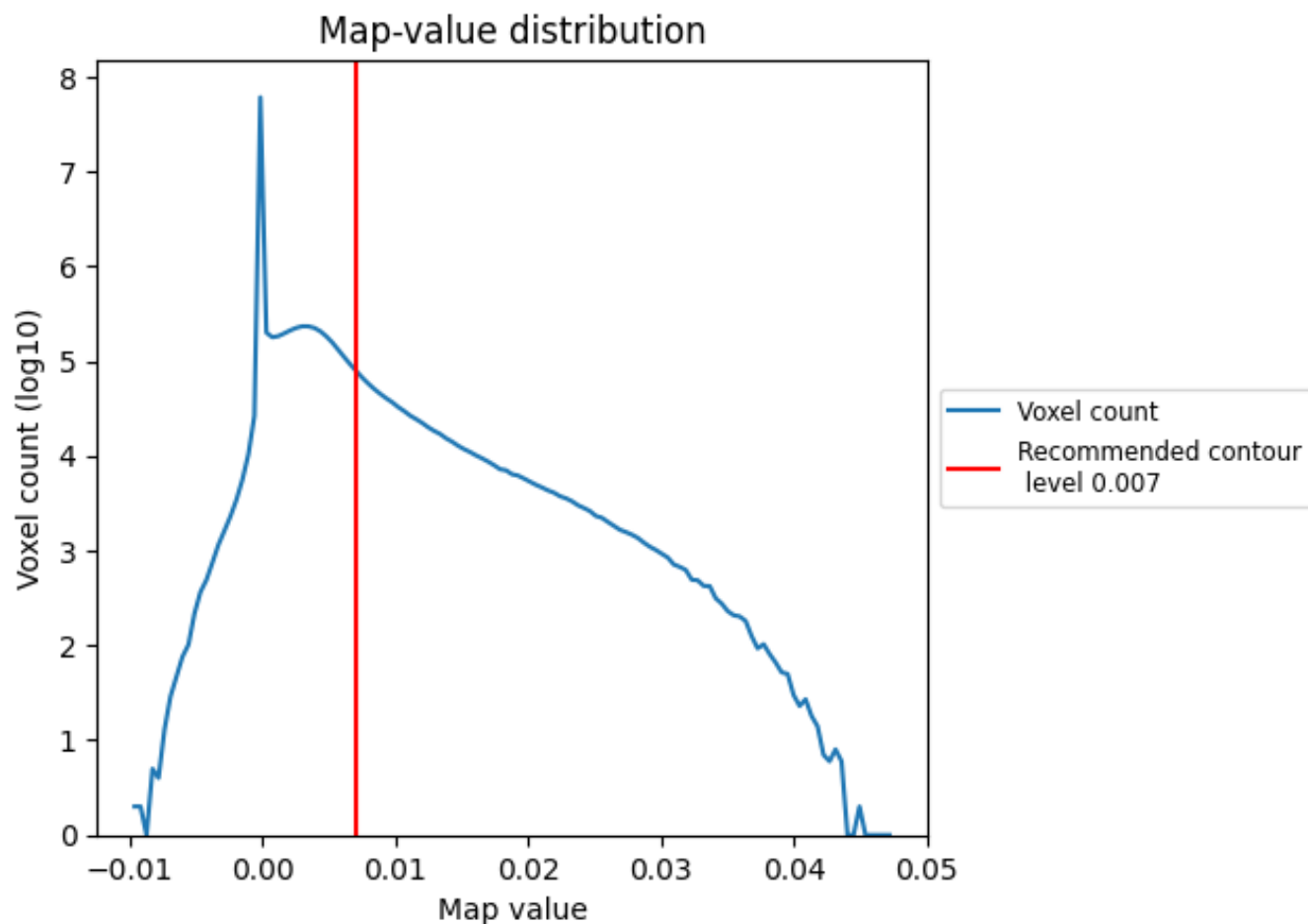
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

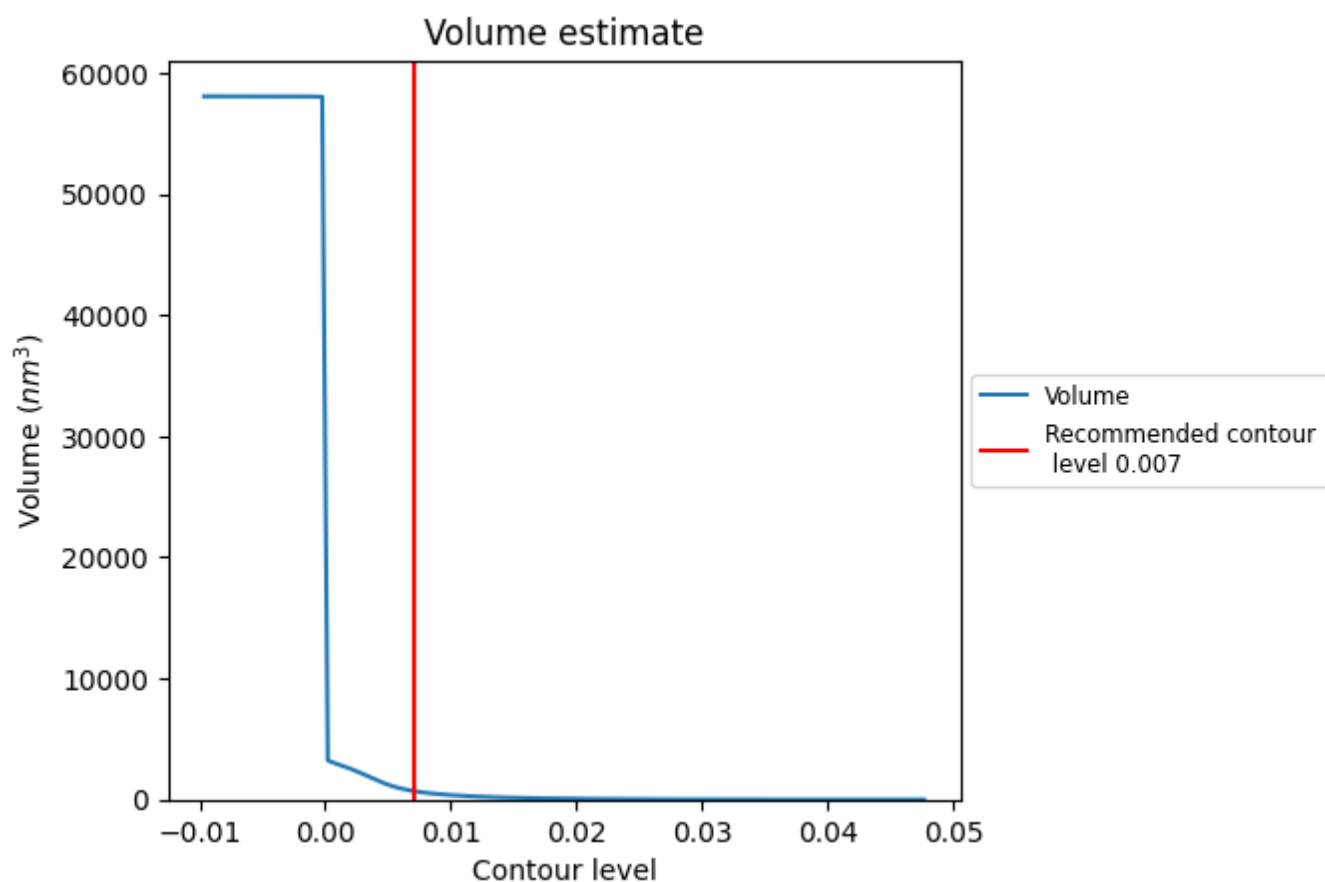
This section contains the results of statistical analysis of the map.

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.

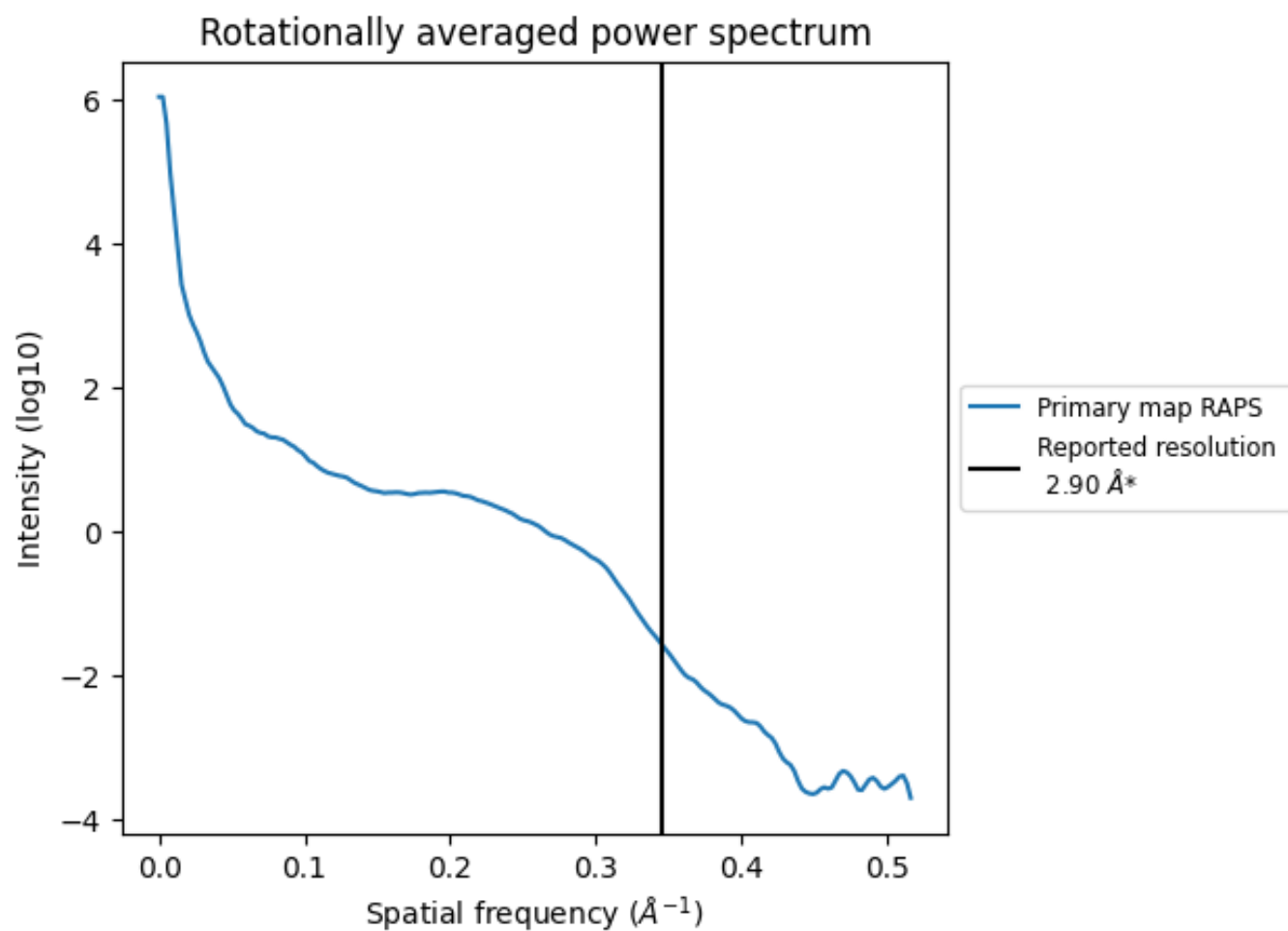
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 706 nm³; this corresponds to an approximate mass of 638 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 0.345 Å⁻¹

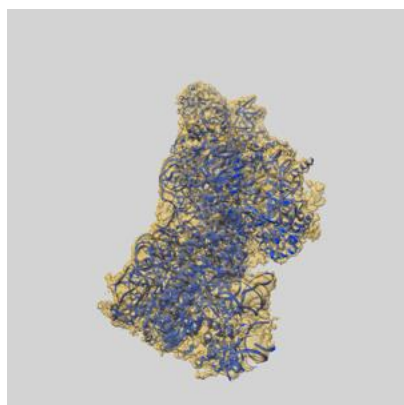
8 Fourier-Shell correlation

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

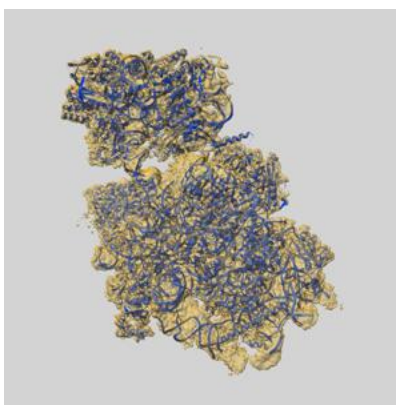
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-16347 and PDB model 8C00. Per-residue inclusion information can be found in section [3](#) on page [8](#).

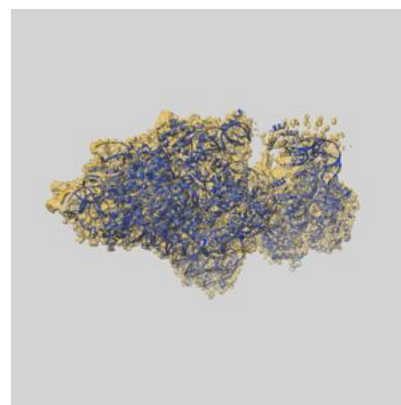
9.1 Map-model overlay [i](#)



X



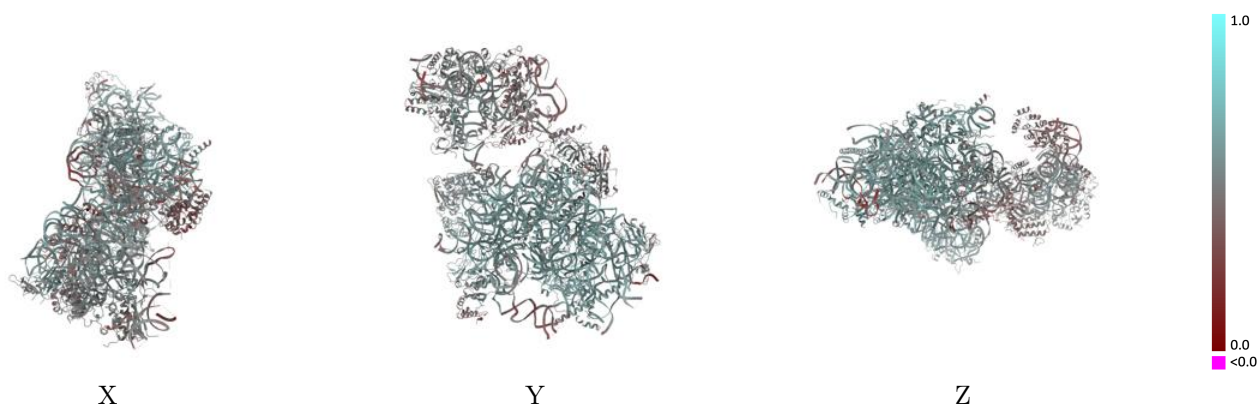
Y



Z

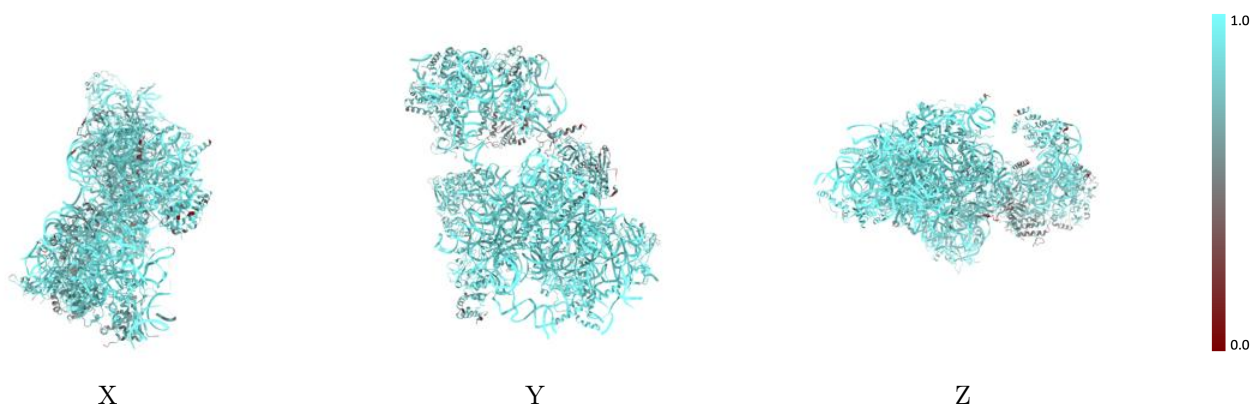
The images above show the 3D surface view of the map at the recommended contour level 0.007 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



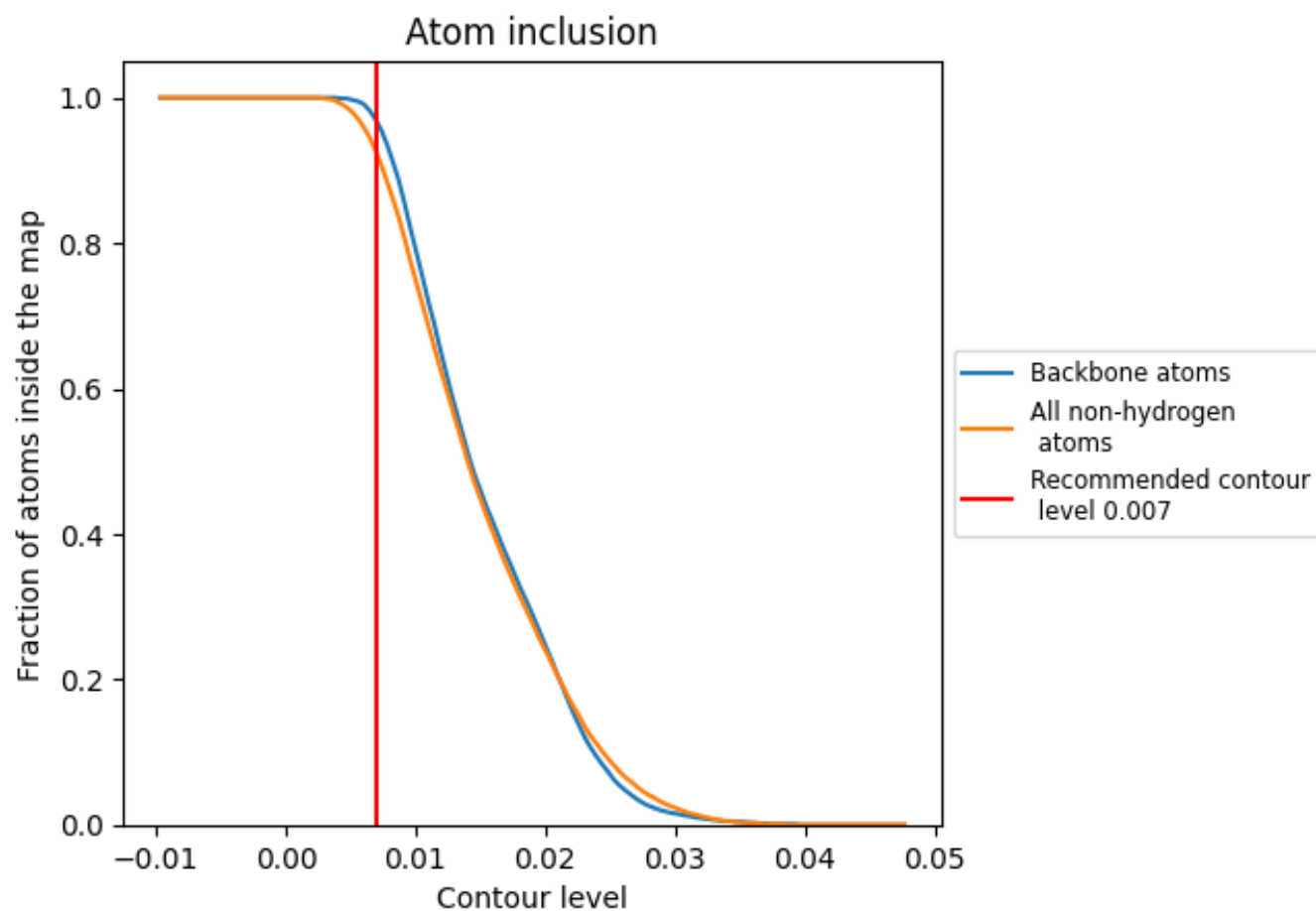
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.007).

9.4 Atom inclusion ⓘ



At the recommended contour level, 97% of all backbone atoms, 92% 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	 0.9240	 0.5350
2	 0.9860	 0.5560
B	 0.8480	 0.4820
E	 0.7600	 0.4720
F	 0.8080	 0.4630
H	 0.7810	 0.4740
I	 0.8560	 0.4940
K	 0.6830	 0.4670
L	 0.8210	 0.4510
Q	 0.9080	 0.5350
S	 0.9720	 0.5970
T	 0.9280	 0.5450
U	 0.8290	 0.5030
V	 0.9690	 0.5940
W	 0.9530	 0.5790
X	 0.9670	 0.5940
Y	 0.9470	 0.5690
Z	 0.9580	 0.5460
b	 0.9740	 0.5900
c	 0.9240	 0.5530
d	 0.9530	 0.5770
e	 0.8410	 0.4080
f	 0.9100	 0.5460
g	 0.9300	 0.5570
p	 0.9040	 0.5130
r	 0.5590	 0.3870
t	 0.7810	 0.4640

