



wwPDB X-ray Structure Validation Summary Report

Mar 1, 2014 – 03:29 AM GMT

PDB ID : 4DR7
Title : Crystal structure of the *Thermus thermophilus* (HB8) 30S ribosomal subunit with codon, crystallographically disordered near-cognate transfer RNA anti-codon stem-loop mismatched at the second codon position, and streptomycin bound
Authors : Demirci, H.; Murphy IV, F.; Murphy, E.; Gregory, S.T.; Dahlberg, A.E.; Jogl, G.
Deposited on : 2012-02-16
Resolution : 3.75 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

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

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

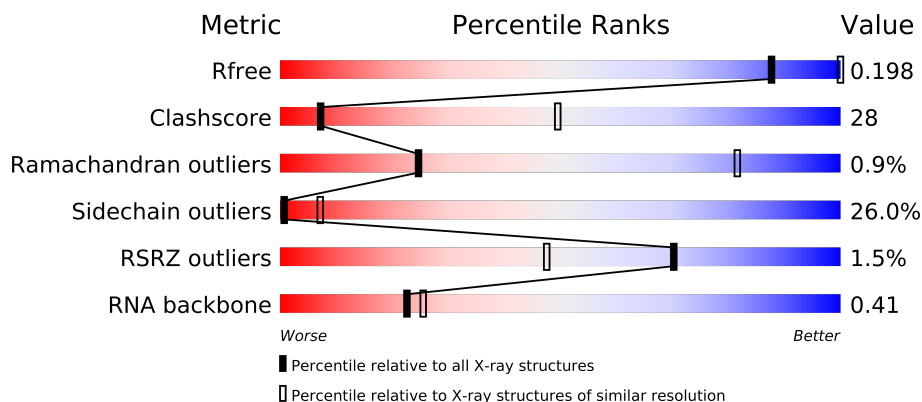
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.75 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.



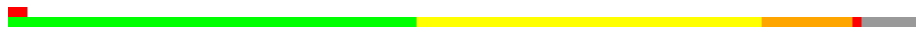
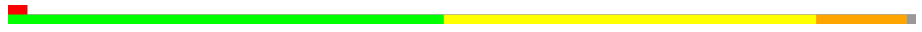

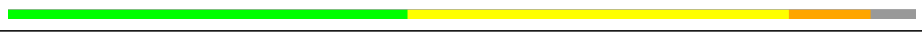
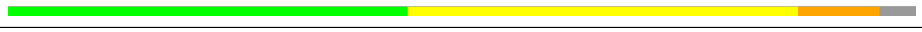
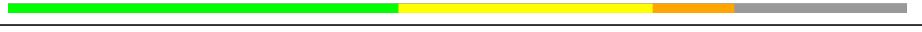
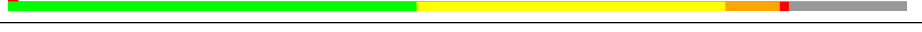



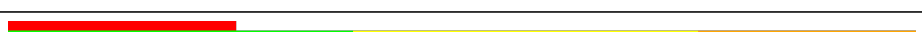
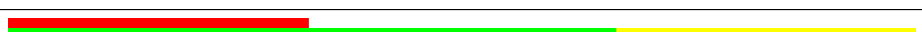

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1134 (4.12-3.40)
Clashscore	79885	1067 (4.02-3.50)
Ramachandran outliers	78287	1017 (4.02-3.50)
Sidechain outliers	78261	1010 (4.02-3.50)
RSRZ outliers	66119	1135 (4.12-3.40)
RNA backbone	1838	1008 (4.52-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	1522	
2	B	256	
3	C	239	
4	D	209	
5	E	162	
6	F	101	
7	G	156	
8	H	138	
9	I	128	
10	J	105	
11	K	129	
12	L	135	

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Mol	Chain	Length	Quality of chain
13	M	126	
14	N	61	
15	O	89	
16	P	88	
17	Q	105	
18	R	88	
19	S	93	
20	T	106	
21	U	27	
22	V	4	
23	W	11	
24	a	8	
25	b	3	

2 Entry composition

There are 29 unique types of molecules in this entry. The entry contains 53659 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 16S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1513	Total	C	N	O	P	0	8	0
			32707	14570	6056	10561	1520			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1534	C	A	CONFLICT	GB M26923.1
A	1535	A	C	CONFLICT	GB M26923.1

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	236	Total	C	N	O	S	0	0	1
			1896	1211	337	343	5			

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	207	Total	C	N	O	S	0	0	1
			1613	1016	315	281	1			

- Molecule 4 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	151	Total	C	N	O	S	0	0	1
			1147	724	218	201	4			

- Molecule 6 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

- Molecule 7 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

- Molecule 8 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	I	127	Total	C	N	O	0	0	0
			1010	639	197	174			

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	99	Total	C	N	O	S	0	0	1
			793	498	157	137	1			

- Molecule 11 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	117	Total	C	N	O	S	0	0	0
			873	543	166	161	3			

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	125	Total	C	N	O	S	0	0	1
			973	612	196	163	2			

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	118	Total	C	N	O	S	0	0	0
			937	579	193	163	2			

- Molecule 14 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

- Molecule 15 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	O	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	P	84	Total	C	N	O	S	0	0	1
			701	443	140	117	1			

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	Q	100	Total	C	N	O	S	0	0	0
			834	534	156	142	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	96	GLN	GLU	CONFLICT	UNP Q5SHP7

- Molecule 18 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	R	71	Total	C	N	O		0	0	0
			585	373	116	96				

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	S	81	Total	C	N	O	S	0	0	1
			648	414	120	112	2			

- Molecule 20 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	T	99	Total	C	N	O	S	0	0	0
			763	470	162	129	2			

- Molecule 21 is a protein called 30S ribosomal protein THX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	U	25	Total	C	N	O	0	0	1
			209	128	51	30			

- Molecule 22 is a RNA chain called 5'-R(*UP*UP*UP*U)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	V	4	Total	C	N	O	P	0	0	0
			77	36	8	30	3			

- Molecule 23 is a RNA chain called 5'-R(*GP*CP*CP*UP*GP*GP*AP*AP*AP*GP*(PSU))-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	W	11	Total	C	N	O	P	0	0	0
			235	106	45	74	10			

- Molecule 24 is a RNA chain called 5'-R(P*UP*GP*GP*AP*AP*AP*GP*(PSU))-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	a	8	Total	C	N	O	P	0	0	0
			175	78	34	55	8			

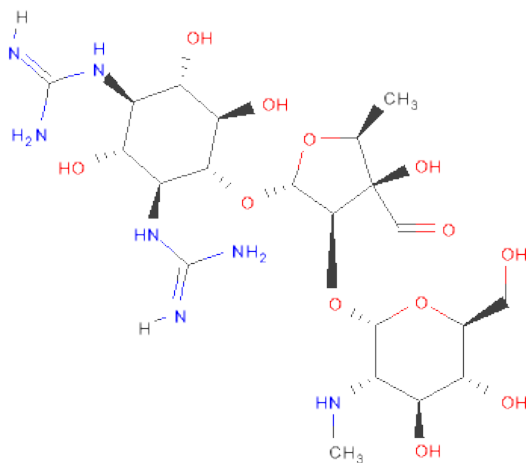
- Molecule 25 is a RNA chain called 5'-R(P*UP*UP*U)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	b	3	Total	C	N	O	P	0	0	0
			60	27	6	24	3			

- Molecule 26 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
26	P	3	Total 3	Mg 3	0	0
26	G	1	Total 1	Mg 1	0	0
26	J	1	Total 1	Mg 1	0	0
26	Q	1	Total 1	Mg 1	0	0
26	D	3	Total 3	Mg 3	0	0
26	E	4	Total 4	Mg 4	0	0
26	H	1	Total 1	Mg 1	0	0
26	A	326	Total 326	Mg 326	0	0
26	N	1	Total 1	Mg 1	0	0
26	S	2	Total 2	Mg 2	0	0
26	F	1	Total 1	Mg 1	0	0

- Molecule 27 is STREPTOMYCIN (three-letter code: SRY) (formula: $C_{21}H_{39}N_7O_{12}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
27	A	1	Total 40	C 21	N 7	O 12	0	0

- Molecule 28 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
28	D	1	Total	Zn	0	0
			1	1		
28	N	1	Total	Zn	0	0
			1	1		

- Molecule 29 is water.

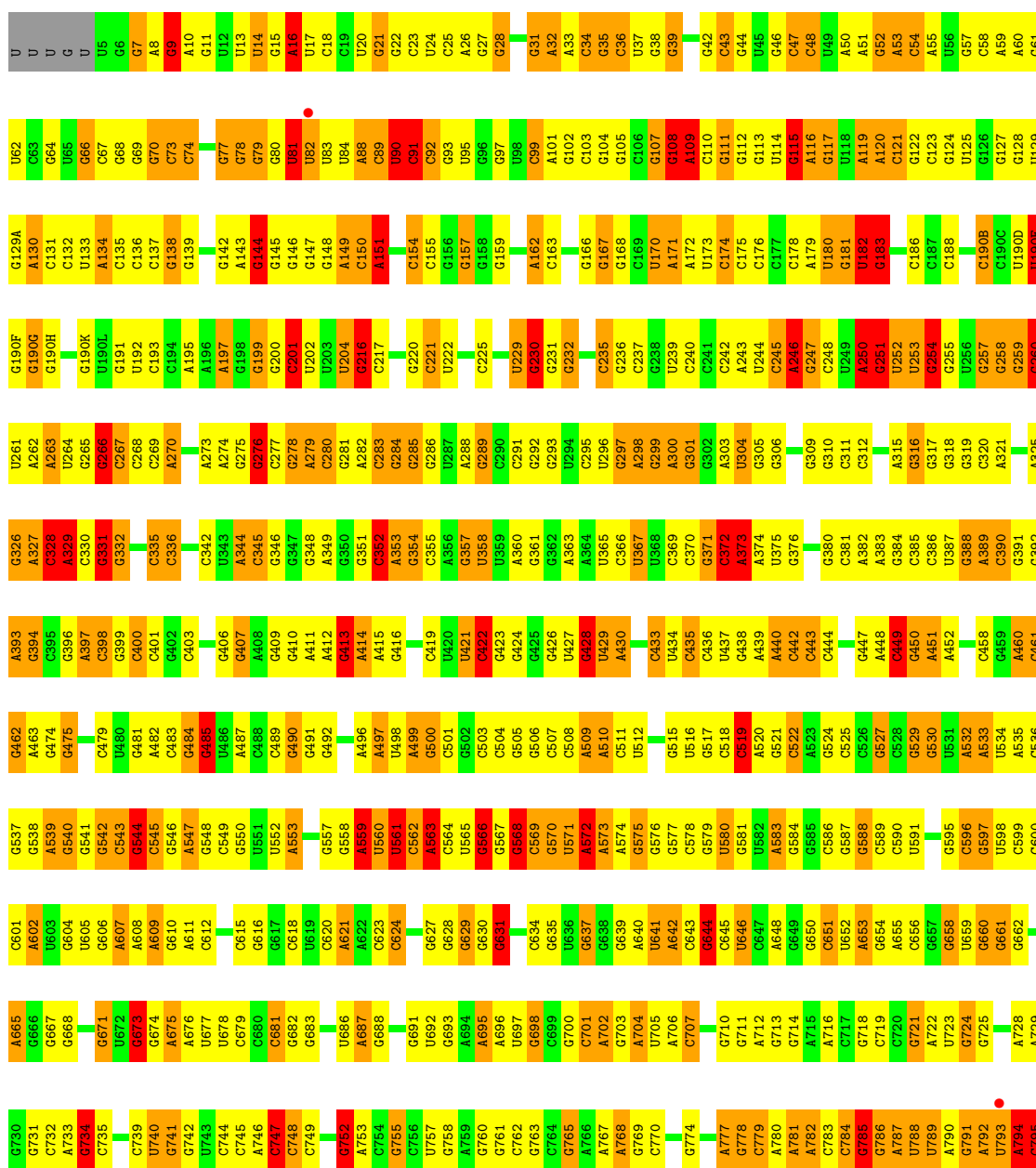
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
29	A	866	Total	O	0	0
			866	866		
29	C	1	Total	O	0	0
			1	1		
29	D	7	Total	O	0	0
			7	7		
29	E	5	Total	O	0	0
			5	5		
29	L	1	Total	O	0	0
			1	1		
29	N	1	Total	O	0	0
			1	1		
29	P	1	Total	O	0	0
			1	1		
29	Q	2	Total	O	0	0
			2	2		
29	T	3	Total	O	0	0
			3	3		
29	U	4	Total	O	0	0
			4	4		
29	W	1	Total	O	0	0
			1	1		

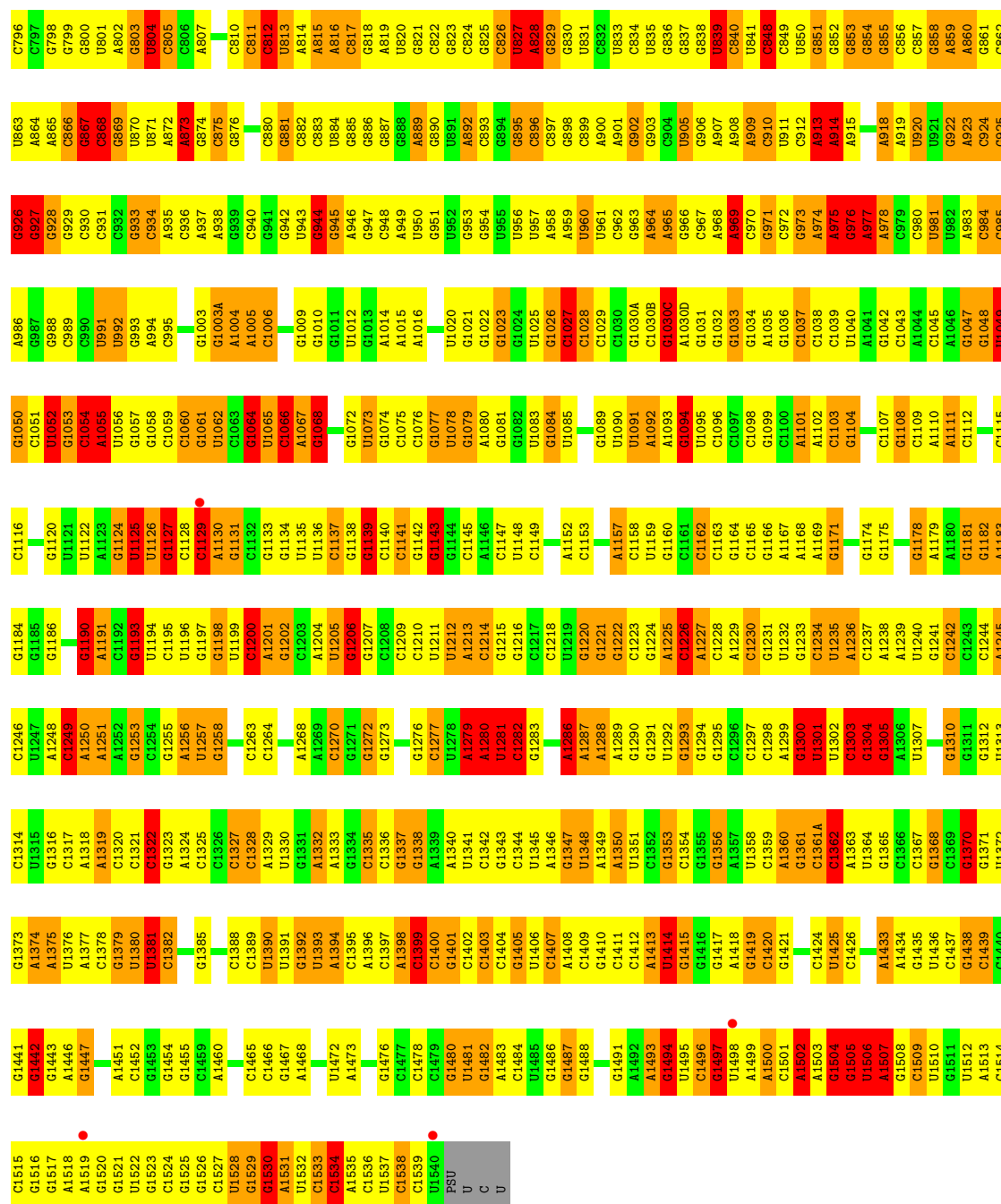
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron 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 dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). 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: 16S rRNA

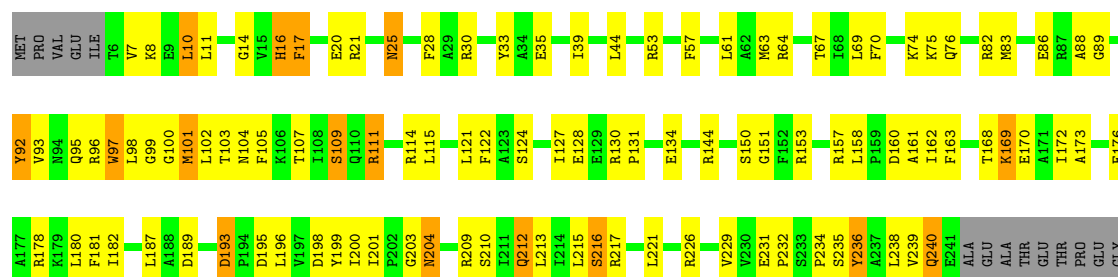
Chain A:





• Molecule 2: 30S ribosomal protein S2

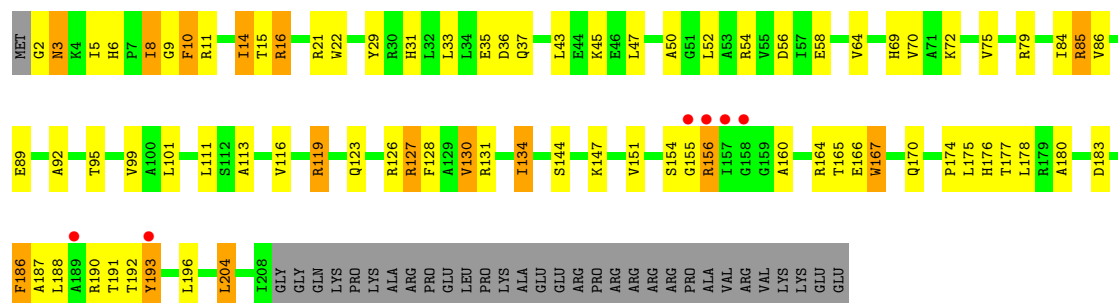
Chain B:



GLU
SER
GLU
VAL
GLU
ALA

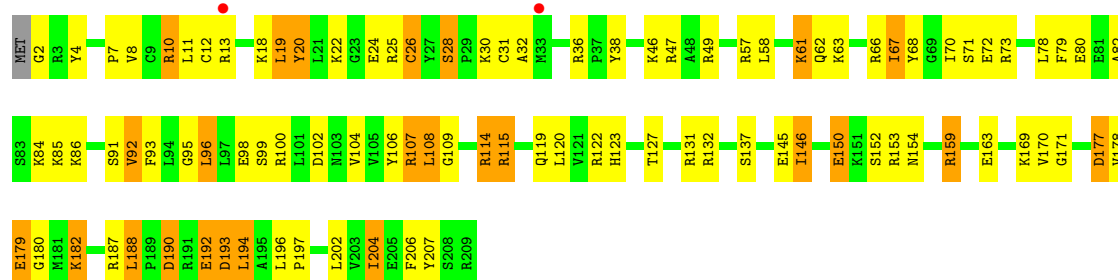
• Molecule 3: 30S ribosomal protein S3

Chain C:



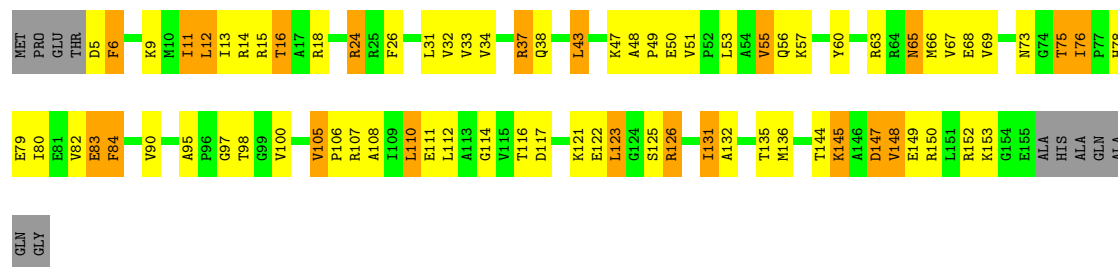
• Molecule 4: 30S ribosomal protein S4

Chain D:



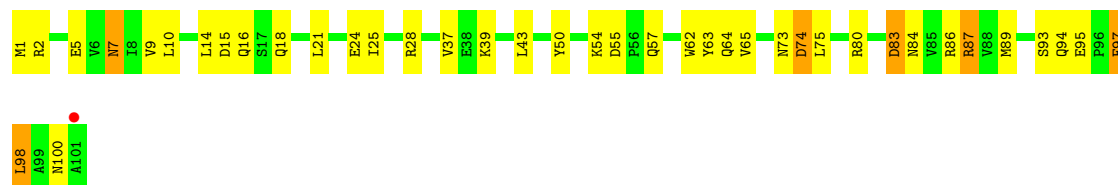
• Molecule 5: 30S ribosomal protein S5

Chain E:



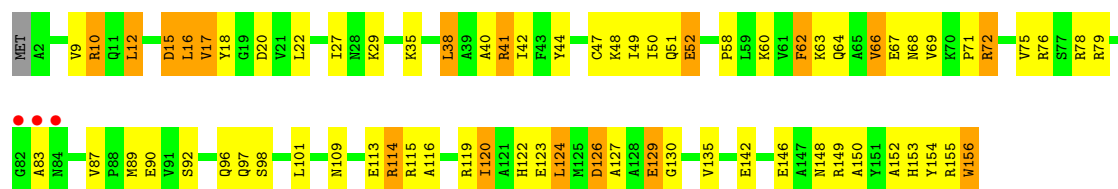
• Molecule 6: 30S ribosomal protein S6

Chain F:



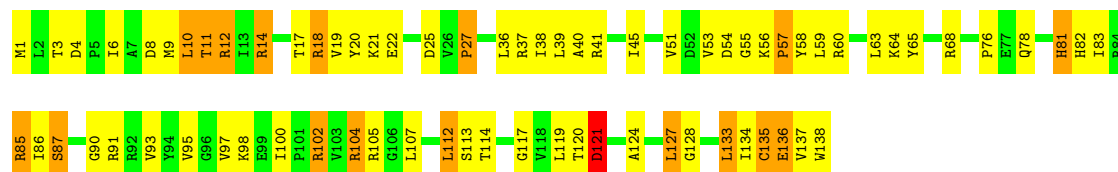
• Molecule 7: 30S ribosomal protein S7

Chain G: 



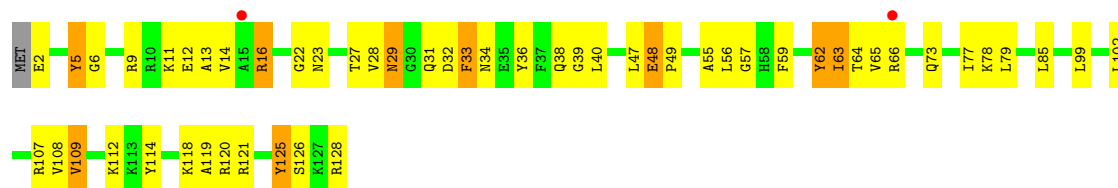
- Molecule 8: 30S ribosomal protein S8

Chain H: 



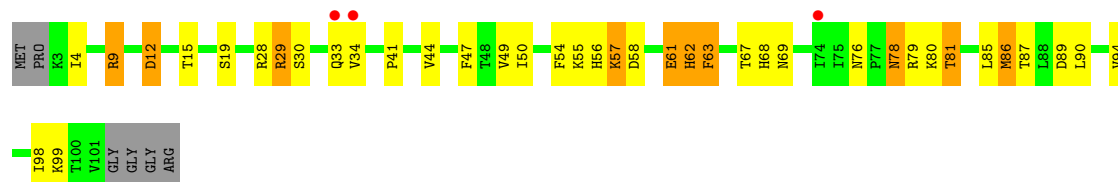
- Molecule 9: 30S ribosomal protein S9

Chain I: 



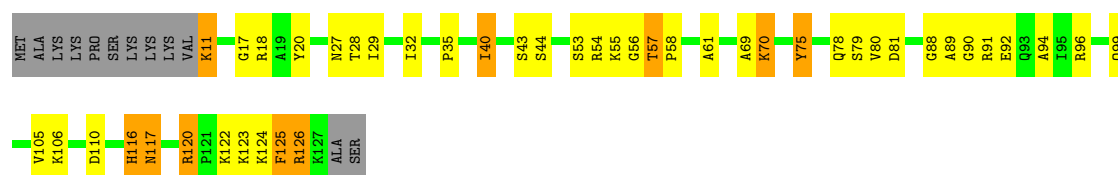
- Molecule 10: 30S ribosomal protein S10

Chain J: 



- Molecule 11: 30S ribosomal protein S11

Chain K: 



- Molecule 12: 30S ribosomal protein S12

Chain L: 





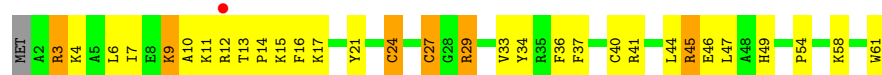
- Molecule 13: 30S ribosomal protein S13

Chain M:



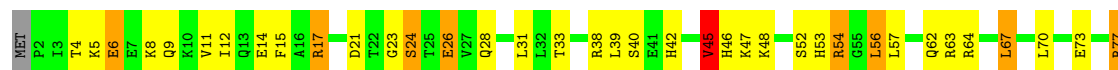
- Molecule 14: 30S ribosomal protein S14

Chain N:



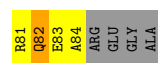
- Molecule 15: 30S ribosomal protein S15

Chain O:



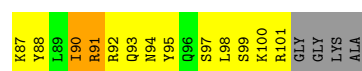
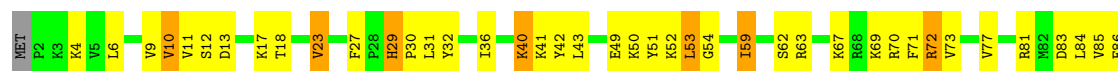
- Molecule 16: 30S ribosomal protein S16

Chain P:



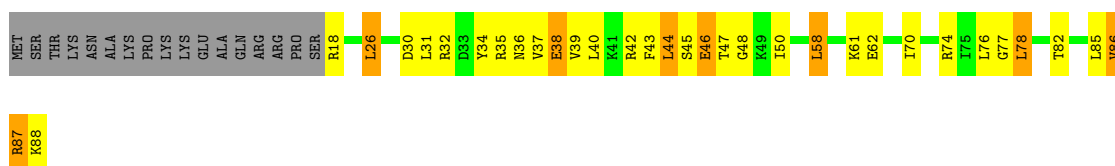
- Molecule 17: 30S ribosomal protein S17

Chain Q:



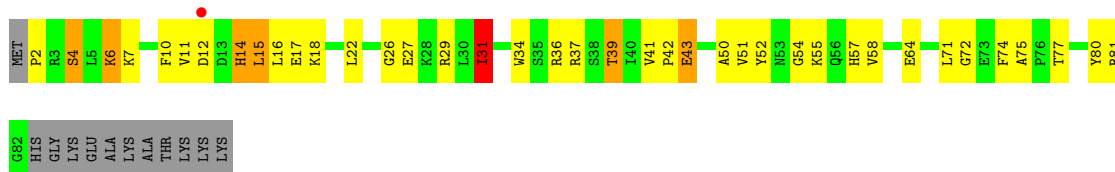
- Molecule 18: 30S ribosomal protein S18

Chain R:



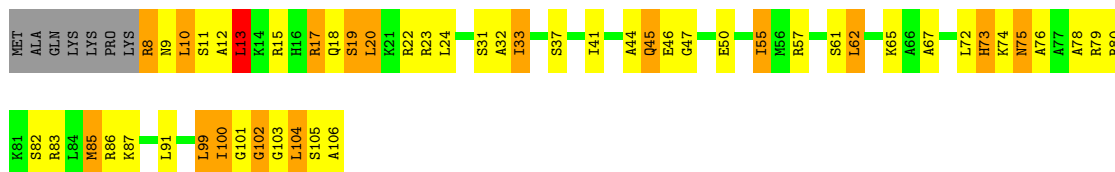
- Molecule 19: 30S ribosomal protein S19

Chain S:



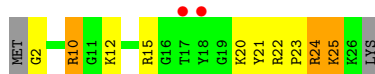
- Molecule 20: 30S ribosomal protein S20

Chain T:



- Molecule 21: 30S ribosomal protein THX

Chain U:



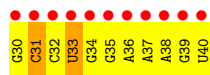
- Molecule 22: 5'-R(*UP*UP*UP*U)-3'

Chain V:



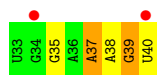
- Molecule 23: 5'-R(*GP*CP*CP*UP*GP*GP*AP*AP*AP*GP*(PSU))-3'

Chain W:



- Molecule 24: 5'-R(P*UP*GP*GP*AP*AP*AP*GP*(PSU))-3'

Chain a:



- Molecule 25: 5'-R(P*UP*UP*U)-3'

Chain b:



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	402.49Å 402.49Å 174.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.91 – 3.75 49.65 – 3.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.91-3.75) 100.0 (49.65-3.75)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 3.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_978)	Depositor
R, R_{free}	0.148 , 0.201 0.146 , 0.198	Depositor DCC
R_{free} test set	7300 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	130.7	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 89.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 146049 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	53659	wwPDB-VP
Average B, all atoms (Å ²)	146.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, M2G, MA6, 0TD, MG, 2MG, 5MC, UR3, 4OC, SRY, 7MG, PSU

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	1.21	135/36234 (0.4%)	1.90	1769/56547 (3.1%)
2	B	0.74	0/1931	0.93	2/2607 (0.1%)
3	C	0.62	0/1637	0.83	0/2207
4	D	0.73	1/1733 (0.1%)	0.93	4/2318 (0.2%)
5	E	1.04	1/1163 (0.1%)	1.17	3/1566 (0.2%)
6	F	0.65	0/856	0.86	0/1154
7	G	0.68	0/1276	0.87	0/1709
8	H	1.11	2/1136 (0.2%)	1.18	4/1527 (0.3%)
9	I	0.65	0/1029	0.88	2/1379 (0.1%)
10	J	0.71	1/806 (0.1%)	0.95	2/1084 (0.2%)
11	K	0.76	0/888	0.97	0/1198
12	L	0.90	0/978	1.08	3/1308 (0.2%)
13	M	0.68	0/947	0.94	0/1270
14	N	0.68	0/501	0.85	1/664 (0.2%)
15	O	0.86	0/745	1.02	3/992 (0.3%)
16	P	0.93	0/717	1.08	3/965 (0.3%)
17	Q	1.08	1/847 (0.1%)	1.25	4/1131 (0.4%)
18	R	0.76	0/590	1.00	1/782 (0.1%)
19	S	0.57	0/662	0.77	0/892
20	T	0.87	0/765	1.18	4/1007 (0.4%)
21	U	0.69	0/213	0.87	0/279
22	V	0.53	0/84	0.98	0/128
23	W	0.62	0/241	0.92	0/375
24	a	0.85	0/174	1.89	10/270 (3.7%)
25	b	0.76	0/65	1.31	2/98 (2.0%)
All	All	1.08	141/56218 (0.3%)	1.66	1817/83457 (2.2%)

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	3
3	C	0	2
4	D	0	1
8	H	0	2
9	I	0	1
10	J	0	2
12	L	0	2
16	P	0	1
19	S	0	1
20	T	0	2
21	U	0	1
All	All	0	18

The worst 5 of 141 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	828	A	N9-C4	-10.24	1.31	1.37
1	A	1513	A	N9-C4	-9.49	1.32	1.37
1	A	266	G	N7-C5	-9.34	1.33	1.39
1	A	573	A	N7-C5	-9.28	1.33	1.39
8	H	135	CYS	CB-SG	-9.04	1.66	1.82

The worst 5 of 1817 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	G	C6-C5-N7	-19.86	118.48	130.40
1	A	366	C	N1-C2-O2	17.84	129.60	118.90
1	A	117	G	N1-C6-O6	15.21	129.03	119.90
1	A	1200	C	C2-N1-C1'	15.12	135.43	118.80
1	A	573	A	C8-N9-C4	-15.02	99.79	105.80

There are no chirality outliers.

5 of 18 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	14	GLY	Peptide
2	B	75	LYS	Peptide
2	B	89	GLY	Peptide
3	C	166	GLU	Peptide
3	C	2	GLY	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	32707	0	0	870	1
2	B	1896	0	6	51	0
3	C	1613	0	0	45	0
4	D	1703	0	0	61	0
5	E	1147	0	0	36	1
6	F	843	0	0	17	0
7	G	1257	0	0	41	0
8	H	1116	0	0	41	0
9	I	1010	0	0	29	0
10	J	793	0	0	25	0
11	K	873	0	12	27	0
12	L	973	0	0	35	0
13	M	937	0	0	37	0
14	N	492	0	0	21	0
15	O	734	0	2	28	0
16	P	701	0	0	22	0
17	Q	834	0	12	32	0
18	R	585	0	12	29	0
19	S	648	0	0	25	0
20	T	763	0	0	31	0
21	U	209	0	0	11	0
22	V	77	0	0	1	0
23	W	235	0	23	23	0
24	a	175	0	0	0	0
25	b	60	0	11	0	0
26	A	326	0	0	0	0
26	D	3	0	0	0	0
26	E	4	0	0	0	0
26	F	1	0	0	0	0
26	G	1	0	0	0	0
26	H	1	0	0	0	0
26	J	1	0	0	0	0
26	N	1	0	0	0	0
26	P	3	0	0	0	0
26	Q	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
26	S	2	0	0	0	0
27	A	40	0	38	12	0
28	D	1	0	0	0	0
28	N	1	0	0	0	0
29	A	866	0	0	91	0
29	C	1	0	0	0	0
29	D	7	0	0	0	0
29	E	5	0	0	0	0
29	L	1	0	0	0	0
29	N	1	0	0	0	0
29	P	1	0	0	0	0
29	Q	2	0	0	0	0
29	T	3	0	0	0	0
29	U	4	0	0	0	0
29	W	1	0	0	0	0
All	All	53659	0	116	1421	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 28.

The worst 5 of 1421 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:451:A:O2'	29:A:2729:HOH:O	1.59	1.20
1:A:692:U:OP1	11:K:124:LYS:NZ	1.82	1.10
1:A:266:G:O2'	1:A:267:C:OP2	1.72	1.08
1:A:328:C:O2'	1:A:329:A:OP2	1.72	1.07
1:A:61:G:O2'	29:A:2002:HOH:O	1.76	1.04

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:82:U:O2	1:A:1400:5MC:N4[3_545]	2.15	0.05
5:E:73:ASN:N	5:E:149:GLU:OE1[7_555]	2.18	0.02

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 X-ray entries followed by that with respect to entries

of similar resolution. 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	234/256 (91%)	197 (84%)	34 (14%)	3 (1%)	18	76
3	C	205/239 (86%)	169 (82%)	35 (17%)	1 (0%)	38	88
4	D	206/209 (99%)	180 (87%)	25 (12%)	1 (0%)	38	88
5	E	149/162 (92%)	137 (92%)	11 (7%)	1 (1%)	30	84
6	F	99/101 (98%)	93 (94%)	6 (6%)	0	100	100
7	G	153/156 (98%)	132 (86%)	21 (14%)	0	100	100
8	H	136/138 (99%)	128 (94%)	7 (5%)	1 (1%)	30	84
9	I	125/128 (98%)	107 (86%)	17 (14%)	1 (1%)	27	83
10	J	97/105 (92%)	77 (79%)	17 (18%)	3 (3%)	7	59
11	K	115/129 (89%)	98 (85%)	17 (15%)	0	100	100
12	L	122/135 (90%)	110 (90%)	8 (7%)	4 (3%)	6	58
13	M	116/126 (92%)	99 (85%)	16 (14%)	1 (1%)	25	81
14	N	58/61 (95%)	50 (86%)	8 (14%)	0	100	100
15	O	86/89 (97%)	72 (84%)	14 (16%)	0	100	100
16	P	82/88 (93%)	74 (90%)	7 (8%)	1 (1%)	19	78
17	Q	98/105 (93%)	93 (95%)	5 (5%)	0	100	100
18	R	69/88 (78%)	60 (87%)	9 (13%)	0	100	100
19	S	79/93 (85%)	68 (86%)	9 (11%)	2 (2%)	9	63
20	T	97/106 (92%)	80 (82%)	16 (16%)	1 (1%)	22	81
21	U	23/27 (85%)	22 (96%)	1 (4%)	0	100	100
All	All	2349/2541 (92%)	2046 (87%)	283 (12%)	20 (1%)	25	81

5 of 20 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	15	THR
9	I	119	ALA
12	L	28	LYS
16	P	83	GLU
19	S	31	ILE

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 X-ray entries followed by that with respect to entries of similar resolution. 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	201/220 (91%)	151 (75%)	50 (25%)	1	8
3	C	160/188 (85%)	119 (74%)	41 (26%)	1	8
4	D	180/181 (99%)	134 (74%)	46 (26%)	1	8
5	E	115/123 (94%)	75 (65%)	40 (35%)	0	3
6	F	90/90 (100%)	60 (67%)	30 (33%)	0	3
7	G	126/127 (99%)	92 (73%)	34 (27%)	1	6
8	H	119/119 (100%)	83 (70%)	36 (30%)	0	5
9	I	98/99 (99%)	75 (76%)	23 (24%)	1	9
10	J	87/92 (95%)	70 (80%)	17 (20%)	2	16
11	K	89/99 (90%)	72 (81%)	17 (19%)	2	16
12	L	103/110 (94%)	80 (78%)	23 (22%)	1	11
13	M	94/101 (93%)	64 (68%)	30 (32%)	0	4
14	N	49/50 (98%)	35 (71%)	14 (29%)	0	5
15	O	79/80 (99%)	60 (76%)	19 (24%)	1	8
16	P	72/74 (97%)	54 (75%)	18 (25%)	1	8
17	Q	95/97 (98%)	74 (78%)	21 (22%)	1	11
18	R	62/77 (80%)	48 (77%)	14 (23%)	1	10
19	S	71/80 (89%)	55 (78%)	16 (22%)	1	11
20	T	76/82 (93%)	51 (67%)	25 (33%)	0	4
21	U	19/22 (86%)	16 (84%)	3 (16%)	4	28
All	All	1985/2111 (94%)	1468 (74%)	517 (26%)	1	7

5 of 517 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
7	G	120	ILE
9	I	63	ILE
19	S	15	LEU
7	G	156	TRP
8	H	95	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1503/1522 (98%)	377 (25%)	48 (3%)
22	V	3/4 (75%)	1 (33%)	0
23	W	10/11 (90%)	2 (20%)	0
24	a	7/8 (87%)	4 (57%)	0
25	b	2/3 (66%)	0	0
All	All	1525/1548 (98%)	384 (25%)	48 (3%)

5 of 384 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	7	G
1	A	9	G
1	A	16	A
1	A	31	G
1	A	32	A

5 of 48 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	587	G
1	A	960	U
1	A	1361(A)	C
1	A	687	A
1	A	812	C

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

18 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	2MG	A	1207	1	24,26,27	2.16	8 (33%)	32,38,41	9.09	2 (6%)
1	5MC	A	1400	1	20,22,23	3.41	6 (30%)	26,32,35	1.47	2 (7%)
1	4OC	A	1402	1	21,23,24	1.78	4 (19%)	26,32,35	1.34	4 (15%)
1	5MC	A	1404	1	20,22,23	2.20	3 (15%)	26,32,35	1.68	4 (15%)
1	5MC	A	1407	1	20,22,23	2.81	4 (20%)	26,32,35	1.33	4 (15%)
1	UR3	A	1498	1	20,22,23	0.94	1 (5%)	23,32,35	0.96	2 (8%)
1	MA6	A	1518[A]	1	26,26,27	1.73	5 (19%)	37,38,41	1.17	4 (10%)
1	MA6	A	1518[B]	1	26,26,27	1.26	2 (7%)	37,38,41	0.98	2 (5%)
1	MA6	A	1519[A]	1	26,26,27	1.12	4 (15%)	37,38,41	1.16	3 (8%)
1	MA6	A	1519[B]	1	26,26,27	1.74	8 (30%)	37,38,41	1.00	3 (8%)
1	PSU	A	1540	1	19,21,22	1.17	1 (5%)	23,30,33	0.94	1 (4%)
1	PSU	A	516	1	19,21,22	1.45	2 (10%)	23,30,33	1.09	3 (13%)
1	7MG	A	527	1	24,26,27	4.46	9 (37%)	34,39,42	1.64	8 (23%)
1	M2G	A	966	1	25,27,28	1.15	3 (12%)	34,40,43	7.33	5 (14%)
1	5MC	A	967	1	20,22,23	1.12	1 (5%)	26,32,35	1.51	4 (15%)
12	0TD	L	92	12	9,9,10	7.57	2 (22%)	9,11,13	3.39	5 (55%)
23	PSU	W	40	23	19,21,22	1.00	1 (5%)	23,30,33	1.22	3 (13%)
24	PSU	a	40	24,1	19,21,22	1.31	1 (5%)	23,30,33	1.36	3 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	2MG	A	1207	1	-	0/10/27/28	0/1/3/3
1	5MC	A	1400	1	-	1/6/25/26	0/2/2/2
1	4OC	A	1402	1	-	2/10/29/30	0/2/2/2
1	5MC	A	1404	1	-	0/6/25/26	0/2/2/2
1	5MC	A	1407	1	-	0/6/25/26	0/2/2/2
1	UR3	A	1498	1	-	0/6/25/26	0/2/2/2
1	MA6	A	1518[A]	1	-	0/13/29/30	0/1/3/3
1	MA6	A	1518[B]	1	-	0/13/29/30	0/1/3/3
1	MA6	A	1519[A]	1	-	0/13/29/30	0/1/3/3
1	MA6	A	1519[B]	1	-	1/13/29/30	0/1/3/3
1	PSU	A	1540	1	-	0/8/25/26	0/2/2/2
1	PSU	A	516	1	-	0/8/25/26	0/2/2/2
1	7MG	A	527	1	-	1/8/37/38	0/1/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	M2G	A	966	1	-	0/12/29/30	0/1/3/3
1	5MC	A	967	1	-	0/6/25/26	0/2/2/2
12	0TD	L	92	12	-	0/10/12/14	0/0/0/0
23	PSU	W	40	23	-	1/8/25/26	0/2/2/2
24	PSU	a	40	24,1	-	1/8/25/26	0/2/2/2

The worst 5 of 65 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	L	92	0TD	O-C	22.20	1.26	1.11
1	A	527	7MG	C8-N9	-17.77	1.32	1.46
1	A	1400	5MC	C2-N1	12.46	1.51	1.38
1	A	1407	5MC	C2-N1	8.99	1.48	1.38
1	A	1404	5MC	C2-N1	8.22	1.47	1.38

The worst 5 of 62 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1207	2MG	C6-C5-N7	-51.19	127.25	134.14
1	A	966	M2G	C6-C5-N7	-42.22	128.46	134.14
12	L	92	0TD	CB-CA-N	-6.73	96.33	109.55
1	A	1404	5MC	C6-N1-C2	6.19	121.67	118.62
1	A	1400	5MC	C2-N3-C4	4.88	119.83	115.41

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1402	4OC	N3-C4-N4-CM4
1	A	1402	4OC	OP2-P-O5'-C5'
1	A	527	7MG	OP2-P-O5'-C5'
1	A	1519[B]	MA6	OP2-P-O5'-C5'
24	a	40	PSU	OP2-P-O5'-C5'

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 347 ligands modelled in this entry, 346 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
27	SRY	A	1928	-	42,42,42	2.22	9 (21%)	63,63,63	2.45	22 (34%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
27	SRY	A	1928	-	-	0/22/87/87	0/3/3/3

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
27	A	1928	SRY	CD1-N31	8.77	1.49	1.33
27	A	1928	SRY	CA1-N11	5.41	1.43	1.33
27	A	1928	SRY	C32-CG2	-4.47	1.47	1.52
27	A	1928	SRY	O53-C53	-3.49	1.35	1.44
27	A	1928	SRY	C23-N23	-3.12	1.42	1.47

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	A	1928	SRY	C43-C33-C23	-7.83	99.32	110.44
27	A	1928	SRY	C12-O42-C42	-5.95	98.71	108.25
27	A	1928	SRY	C61-C11-N11	-5.68	98.51	110.56
27	A	1928	SRY	C13-O13-C22	-5.46	106.79	116.28
27	A	1928	SRY	C42-C32-C22	-4.77	96.21	101.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1512/1522 (99%)	-0.41	6 (0%) 90 79	80, 129, 232, 327	0
2	B	236/256 (92%)	0.14	0 100 100	62, 153, 220, 247	0
3	C	207/239 (86%)	0.09	6 (2%) 49 31	94, 191, 226, 244	0
4	D	208/209 (99%)	0.08	2 (0%) 79 58	105, 149, 198, 214	0
5	E	151/162 (93%)	-0.13	0 100 100	80, 114, 166, 193	0
6	F	101/101 (100%)	-0.17	1 (0%) 79 58	110, 150, 182, 226	0
7	G	155/156 (99%)	-0.09	3 (1%) 64 42	121, 160, 223, 238	0
8	H	138/138 (100%)	-0.04	0 100 100	75, 104, 138, 155	0
9	I	127/128 (99%)	0.32	2 (1%) 68 46	125, 184, 217, 240	0
10	J	99/105 (94%)	0.37	3 (3%) 48 30	81, 203, 272, 295	0
11	K	117/129 (90%)	-0.17	0 100 100	88, 130, 160, 167	0
12	L	124/135 (91%)	0.35	5 (4%) 36 24	97, 136, 168, 247	0
13	M	118/126 (93%)	0.29	2 (1%) 67 45	127, 158, 192, 211	0
14	N	60/61 (98%)	0.09	1 (1%) 67 45	144, 169, 210, 246	0
15	O	88/89 (98%)	0.33	2 (2%) 57 37	94, 127, 182, 225	0
16	P	84/88 (95%)	0.12	0 100 100	96, 126, 164, 239	0
17	Q	100/105 (95%)	0.09	0 100 100	83, 111, 150, 189	0
18	R	71/88 (80%)	0.04	0 100 100	95, 132, 178, 229	0
19	S	81/93 (87%)	0.41	1 (1%) 75 53	84, 184, 228, 234	0
20	T	99/106 (93%)	0.10	0 100 100	95, 126, 167, 212	0
21	U	25/27 (92%)	0.72	2 (8%) 12 11	77, 163, 192, 220	0
22	V	4/4 (100%)	3.70	4 (100%) 0 0	263, 267, 272, 275	0
23	W	11/11 (100%)	6.31	11 (100%) 0 0	234, 291, 345, 387	5 (45%)
24	a	8/8 (100%)	1.05	2 (25%) 1 2	200, 218, 290, 330	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
25	b	3/3 (100%)	1.25	1 (33%) 1 1	172, 172, 206, 218	0
All	All	3927/4089 (96%)	-0.07	54 (1%) 70 49	62, 140, 222, 387	5 (0%)

The worst 5 of 54 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
23	W	32	C	10.9
1	A	1498	UR3	9.1
23	W	31	C	8.4
23	W	38	A	7.2
23	W	33	U	7.2

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	5MC	A	967	21/22	0.16	-	117,131,145,146	0
1	5MC	A	1404	21/22	0.17	-	102,129,148,149	0
1	7MG	A	527	24/25	0.18	-	91,114,123,126	0
1	M2G	A	966	25/26	0.21	-	122,137,142,145	0
1	2MG	A	1207	24/25	0.14	-	154,167,200,202	0
1	UR3	A	1498	21/22	0.27	-	111,124,183,193	0
1	MA6	A	1518[A]	24/25	0.20	-	110,122,127,131	24
1	5MC	A	1407	21/22	0.12	-	127,152,158,162	0
12	0TD	L	92	10/11	0.19	-	113,121,127,289	0
23	PSU	W	40	20/21	0.39	-	291,301,325,326	0
1	PSU	A	516	20/21	0.10	-	123,147,168,168	0
1	4OC	A	1402	22/23	0.21	-	104,119,127,142	0
1	PSU	A	1540	20/21	0.35	-	253,269,289,293	0
1	MA6	A	1519[A]	24/25	0.29	-	100,115,125,126	24
1	MA6	A	1518[B]	24/25	0.20	-	107,122,137,148	24
1	MA6	A	1519[B]	24/25	0.29	-	101,116,129,130	24
24	PSU	a	40	20/21	0.37	-	208,236,258,262	0
1	5MC	A	1400	21/22	0.20	-	103,130,148,159	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
26	MG	A	1770	1/1	0.51	-	106,106,106,106	0
26	MG	A	1862	1/1	0.30	-	268,268,268,268	0
26	MG	A	1617	1/1	0.23	-	85,85,85,85	0
26	MG	A	1925	1/1	0.16	-	114,114,114,114	0
26	MG	A	1917	1/1	0.77	-	128,128,128,128	0
26	MG	A	1644	1/1	0.34	-	127,127,127,127	0
26	MG	A	1623	1/1	0.12	-	67,67,67,67	0
26	MG	A	1755	1/1	0.48	-	74,74,74,74	0
26	MG	A	1815	1/1	0.12	-	506,506,506,506	0
26	MG	A	1885	1/1	0.15	-	84,84,84,84	0
26	MG	A	1800	1/1	0.19	-	415,415,415,415	1
26	MG	P	103	1/1	0.49	-	96,96,96,96	0
26	MG	A	1859	1/1	0.17	-	427,427,427,427	0
28	ZN	N	101	1/1	0.15	-	164,164,164,164	0
26	MG	A	1871	1/1	0.24	-	392,392,392,392	0
26	MG	A	1832	1/1	0.09	-	278,278,278,278	0
26	MG	A	1669	1/1	0.12	-	123,123,123,123	0
26	MG	A	1905	1/1	0.24	-	105,105,105,105	0
26	MG	A	1756	1/1	0.08	-	126,126,126,126	0
26	MG	E	203	1/1	0.12	-	101,101,101,101	0
26	MG	A	1716	1/1	0.23	-	132,132,132,132	0
26	MG	A	1845	1/1	0.21	-	410,410,410,410	0
26	MG	A	1604	1/1	0.30	-	91,91,91,91	0
26	MG	A	1926	1/1	0.56	-	116,116,116,116	0
26	MG	A	1873	1/1	0.05	-	361,361,361,361	0
26	MG	A	1622	1/1	0.66	-	81,81,81,81	0
26	MG	A	1609	1/1	0.07	-	157,157,157,157	0
26	MG	A	1771	1/1	0.50	-	102,102,102,102	0
26	MG	A	1918	1/1	0.19	-	75,75,75,75	0
26	MG	A	1676	1/1	0.34	-	158,158,158,158	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1731	1/1	0.19	-	68,68,68,68	0
26	MG	A	1745	1/1	0.56	-	92,92,92,92	0
26	MG	A	1662	1/1	0.13	-	123,123,123,123	0
26	MG	A	1906	1/1	0.12	-	64,64,64,64	0
26	MG	A	1680	1/1	0.45	-	214,214,214,214	0
26	MG	A	1822	1/1	0.12	-	293,293,293,293	0
26	MG	A	1757	1/1	0.24	-	88,88,88,88	0
26	MG	A	1659	1/1	0.31	-	90,90,90,90	0
26	MG	A	1834	1/1	0.84	-	412,412,412,412	1
26	MG	A	1827	1/1	0.09	-	251,251,251,251	0
26	MG	A	1851	1/1	0.29	-	328,328,328,328	0
26	MG	A	1694	1/1	0.49	-	87,87,87,87	0
26	MG	A	1908	1/1	0.34	-	81,81,81,81	0
26	MG	A	1899	1/1	0.40	-	68,68,68,68	0
26	MG	A	1639	1/1	0.21	-	87,87,87,87	0
26	MG	A	1700	1/1	0.38	-	302,302,302,302	0
26	MG	A	1717	1/1	0.30	-	152,152,152,152	0
26	MG	A	1738	1/1	0.05	-	68,68,68,68	0
26	MG	A	1865	1/1	0.34	-	457,457,457,457	0
26	MG	A	1894	1/1	0.24	-	94,94,94,94	0
26	MG	A	1737	1/1	0.71	-	89,89,89,89	0
26	MG	A	1740	1/1	0.66	-	93,93,93,93	0
26	MG	A	1900	1/1	0.39	-	93,93,93,93	0
26	MG	A	1902	1/1	0.39	-	119,119,119,119	0
26	MG	A	1627	1/1	0.16	-	141,141,141,141	0
26	MG	A	1817	1/1	0.18	-	419,419,419,419	0
26	MG	A	1811	1/1	1.17	-	468,468,468,468	0
26	MG	A	1730	1/1	0.65	-	97,97,97,97	0
26	MG	A	1723	1/1	0.29	-	117,117,117,117	0
26	MG	A	1765	1/1	0.17	-	104,104,104,104	0
26	MG	A	1849	1/1	0.26	-	471,471,471,471	0
26	MG	A	1882	1/1	0.33	-	315,315,315,315	0
26	MG	A	1701	1/1	0.21	-	108,108,108,108	0
26	MG	A	1857	1/1	0.29	-	308,308,308,308	0
26	MG	A	1883	1/1	0.33	-	443,443,443,443	0
26	MG	A	1741	1/1	0.16	-	105,105,105,105	0
26	MG	A	1689	1/1	0.12	-	126,126,126,126	0
26	MG	A	1613	1/1	0.13	-	202,202,202,202	0
26	MG	A	1664	1/1	0.26	-	154,154,154,154	0
26	MG	A	1711	1/1	0.86	-	133,133,133,133	0
26	MG	A	1624	1/1	1.09	-	147,147,147,147	0
26	MG	Q	201	1/1	0.57	-	145,145,145,145	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
26	MG	A	1615	1/1	0.11	-	147,147,147,147	0
26	MG	A	1843	1/1	0.49	-	357,357,357,357	0
26	MG	A	1751	1/1	0.54	-	76,76,76,76	0
26	MG	A	1797	1/1	0.12	-	502,502,502,502	0
26	MG	A	1656	1/1	0.26	-	236,236,236,236	0
26	MG	A	1660	1/1	0.13	-	224,224,224,224	0
26	MG	A	1821	1/1	0.21	-	428,428,428,428	0
26	MG	A	1835	1/1	0.20	-	309,309,309,309	0
28	ZN	D	301	1/1	0.28	-	138,138,138,138	0
26	MG	A	1667	1/1	0.19	-	119,119,119,119	0
26	MG	G	201	1/1	0.84	-	117,117,117,117	0
26	MG	A	1640	1/1	0.26	-	172,172,172,172	0
26	MG	A	1907	1/1	0.32	-	120,120,120,120	0
26	MG	A	1632	1/1	0.38	-	248,248,248,248	0
26	MG	A	1606	1/1	0.40	-	87,87,87,87	0
26	MG	N	102	1/1	0.64	-	107,107,107,107	0
26	MG	A	1781	1/1	0.34	-	104,104,104,104	0
26	MG	A	1796	1/1	0.12	-	416,416,416,416	0
26	MG	A	1829	1/1	0.29	-	489,489,489,489	0
26	MG	A	1921	1/1	0.41	-	84,84,84,84	0
26	MG	A	1657	1/1	0.27	-	220,220,220,220	0
26	MG	A	1603	1/1	0.14	-	277,277,277,277	0
26	MG	A	1896	1/1	0.17	-	114,114,114,114	0
26	MG	A	1649	1/1	0.34	-	84,84,84,84	0
26	MG	A	1725	1/1	0.26	-	92,92,92,92	0
26	MG	A	1682	1/1	0.07	-	242,242,242,242	0
26	MG	A	1888	1/1	0.27	-	89,89,89,89	0
26	MG	A	1807	1/1	0.47	-	517,517,517,517	1
26	MG	A	1860	1/1	0.16	-	460,460,460,460	0
26	MG	A	1626	1/1	0.23	-	245,245,245,245	0
26	MG	A	1668	1/1	0.28	-	173,173,173,173	0
26	MG	A	1855	1/1	0.13	-	401,401,401,401	0
26	MG	A	1612	1/1	0.34	-	281,281,281,281	0
26	MG	A	1816	1/1	0.41	-	454,454,454,454	0
26	MG	A	1853	1/1	0.27	-	502,502,502,502	0
26	MG	A	1812	1/1	0.57	-	444,444,444,444	0
26	MG	A	1858	1/1	0.23	-	355,355,355,355	0
26	MG	A	1641	1/1	0.18	-	102,102,102,102	0
26	MG	A	1633	1/1	0.22	-	118,118,118,118	0
26	MG	A	1675	1/1	1.13	-	122,122,122,122	0
26	MG	A	1690	1/1	0.08	-	114,114,114,114	0
26	MG	A	1753	1/1	0.31	-	109,109,109,109	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1870	1/1	0.35	-	406,406,406,406	0
26	MG	H	201	1/1	0.36	-	67,67,67,67	0
26	MG	A	1743	1/1	0.44	-	105,105,105,105	0
26	MG	A	1780	1/1	0.31	-	119,119,119,119	0
26	MG	A	1910	1/1	0.29	-	120,120,120,120	0
26	MG	A	1914	1/1	1.00	-	99,99,99,99	0
26	MG	A	1608	1/1	0.24	-	73,73,73,73	0
26	MG	A	1650	1/1	0.18	-	151,151,151,151	0
26	MG	A	1848	1/1	0.29	-	407,407,407,407	0
26	MG	A	1763	1/1	0.18	-	120,120,120,120	0
26	MG	A	1801	1/1	0.17	-	464,464,464,464	0
26	MG	A	1838	1/1	0.57	-	538,538,538,538	0
26	MG	A	1688	1/1	0.53	-	186,186,186,186	0
26	MG	A	1824	1/1	0.14	-	503,503,503,503	0
26	MG	A	1788	1/1	0.12	-	285,285,285,285	0
26	MG	A	1864	1/1	0.18	-	393,393,393,393	0
26	MG	A	1878	1/1	0.17	-	346,346,346,346	0
26	MG	A	1779	1/1	0.27	-	102,102,102,102	0
26	MG	A	1830	1/1	0.17	-	480,480,480,480	0
26	MG	A	1729	1/1	0.88	-	101,101,101,101	0
26	MG	A	1791	1/1	0.17	-	283,283,283,283	0
26	MG	A	1844	1/1	0.07	-	420,420,420,420	0
26	MG	A	1739	1/1	0.12	-	78,78,78,78	0
26	MG	A	1746	1/1	0.15	-	104,104,104,104	0
26	MG	A	1880	1/1	0.45	-	508,508,508,508	0
26	MG	A	1708	1/1	0.24	-	152,152,152,152	0
26	MG	A	1715	1/1	0.21	-	229,229,229,229	0
26	MG	A	1724	1/1	0.62	-	117,117,117,117	0
26	MG	A	1911	1/1	0.10	-	100,100,100,100	0
26	MG	A	1607	1/1	0.15	-	154,154,154,154	0
26	MG	A	1786	1/1	0.06	-	109,109,109,109	0
26	MG	A	1634	1/1	0.10	-	107,107,107,107	0
26	MG	A	1679	1/1	0.14	-	369,369,369,369	0
26	MG	A	1698	1/1	0.20	-	124,124,124,124	0
26	MG	A	1666	1/1	0.09	-	149,149,149,149	0
26	MG	A	1852	1/1	0.31	-	450,450,450,450	0
26	MG	A	1794	1/1	0.24	-	518,518,518,518	0
26	MG	A	1828	1/1	0.15	-	432,432,432,432	0
26	MG	A	1923	1/1	0.81	-	86,86,86,86	0
26	MG	A	1727	1/1	0.61	-	84,84,84,84	0
26	MG	A	1768	1/1	0.40	-	107,107,107,107	0
26	MG	A	1819	1/1	0.46	-	493,493,493,493	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1793	1/1	0.14	-	483,483,483,483	0
26	MG	A	1774	1/1	0.52	-	109,109,109,109	0
26	MG	A	1890	1/1	0.14	-	112,112,112,112	0
26	MG	A	1814	1/1	0.10	-	440,440,440,440	0
26	MG	A	1803	1/1	0.17	-	345,345,345,345	0
26	MG	A	1776	1/1	0.12	-	96,96,96,96	0
26	MG	A	1920	1/1	0.08	-	138,138,138,138	0
26	MG	A	1671	1/1	0.24	-	96,96,96,96	0
26	MG	A	1636	1/1	0.52	-	88,88,88,88	0
26	MG	A	1898	1/1	0.53	-	89,89,89,89	0
26	MG	A	1887	1/1	0.48	-	102,102,102,102	0
26	MG	A	1837	1/1	0.15	-	469,469,469,469	0
26	MG	A	1823	1/1	0.10	-	422,422,422,422	0
26	MG	A	1658	1/1	0.18	-	115,115,115,115	0
26	MG	A	1904	1/1	0.75	-	97,97,97,97	0
26	MG	A	1642	1/1	0.29	-	78,78,78,78	0
26	MG	S	101	1/1	0.12	-	115,115,115,115	0
26	MG	A	1621	1/1	0.20	-	131,131,131,131	0
26	MG	A	1678	1/1	0.11	-	217,217,217,217	0
26	MG	A	1736	1/1	0.15	-	123,123,123,123	0
26	MG	A	1638	1/1	0.12	-	86,86,86,86	0
26	MG	A	1866	1/1	0.20	-	434,434,434,434	0
26	MG	A	1901	1/1	0.31	-	101,101,101,101	0
26	MG	A	1661	1/1	0.21	-	244,244,244,244	0
26	MG	A	1804	1/1	0.38	-	447,447,447,447	0
26	MG	A	1919	1/1	0.20	-	78,78,78,78	0
26	MG	A	1767	1/1	0.28	-	111,111,111,111	0
26	MG	A	1783	1/1	0.12	-	110,110,110,110	0
26	MG	A	1663	1/1	0.18	-	160,160,160,160	0
26	MG	A	1686	1/1	0.28	-	184,184,184,184	0
26	MG	A	1778	1/1	0.10	-	119,119,119,119	1
26	MG	J	201	1/1	0.34	-	109,109,109,109	0
26	MG	A	1839	1/1	0.08	-	467,467,467,467	0
26	MG	A	1712	1/1	0.19	-	170,170,170,170	0
26	MG	A	1808	1/1	0.56	-	498,498,498,498	0
26	MG	A	1775	1/1	0.10	-	79,79,79,79	0
26	MG	A	1710	1/1	0.13	-	253,253,253,253	0
26	MG	A	1630	1/1	0.26	-	158,158,158,158	0
26	MG	A	1876	1/1	0.37	-	463,463,463,463	1
26	MG	A	1720	1/1	0.29	-	74,74,74,74	0
26	MG	A	1881	1/1	0.13	-	414,414,414,414	0
26	MG	A	1696	1/1	0.25	-	263,263,263,263	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1611	1/1	0.09	-	113,113,113,113	0
26	MG	F	601	1/1	0.07	-	102,102,102,102	0
26	MG	A	1785	1/1	0.14	-	108,108,108,108	0
26	MG	A	1836	1/1	0.21	-	416,416,416,416	1
26	MG	A	1893	1/1	0.69	-	114,114,114,114	0
26	MG	A	1648	1/1	0.28	-	93,93,93,93	0
26	MG	P	101	1/1	0.35	-	58,58,58,58	0
26	MG	A	1699	1/1	0.08	-	414,414,414,414	0
26	MG	A	1818	1/1	0.12	-	189,189,189,189	0
26	MG	A	1895	1/1	0.12	-	106,106,106,106	0
26	MG	A	1748	1/1	0.23	-	110,110,110,110	0
26	MG	A	1750	1/1	0.11	-	74,74,74,74	0
26	MG	A	1897	1/1	0.49	-	114,114,114,114	0
26	MG	A	1916	1/1	0.14	-	119,119,119,119	0
26	MG	A	1654	1/1	0.09	-	161,161,161,161	0
26	MG	A	1773	1/1	0.10	-	103,103,103,103	0
26	MG	A	1787	1/1	0.14	-	496,496,496,496	0
26	MG	A	1820	1/1	0.14	-	373,373,373,373	0
26	MG	E	201	1/1	0.42	-	95,95,95,95	0
26	MG	A	1868	1/1	0.42	-	380,380,380,380	0
26	MG	A	1702	1/1	0.13	-	279,279,279,279	0
26	MG	A	1637	1/1	0.27	-	228,228,228,228	0
26	MG	A	1628	1/1	0.33	-	152,152,152,152	0
26	MG	A	1806	1/1	0.20	-	505,505,505,505	0
26	MG	A	1915	1/1	0.57	-	117,117,117,117	0
26	MG	A	1863	1/1	0.21	-	409,409,409,409	0
26	MG	A	1912	1/1	0.67	-	104,104,104,104	0
26	MG	A	1616	1/1	0.18	-	63,63,63,63	0
26	MG	A	1693	1/1	0.10	-	108,108,108,108	0
26	MG	A	1742	1/1	0.44	-	97,97,97,97	0
26	MG	A	1685	1/1	0.49	-	126,126,126,126	0
26	MG	E	204	1/1	0.22	-	128,128,128,128	0
26	MG	A	1674	1/1	0.04	-	260,260,260,260	0
26	MG	A	1728	1/1	0.97	-	114,114,114,114	0
26	MG	A	1872	1/1	0.25	-	359,359,359,359	0
26	MG	A	1869	1/1	0.20	-	431,431,431,431	0
26	MG	A	1809	1/1	0.22	-	494,494,494,494	0
26	MG	A	1846	1/1	0.21	-	449,449,449,449	0
26	MG	E	202	1/1	0.12	-	124,124,124,124	0
26	MG	A	1605	1/1	0.09	-	254,254,254,254	0
26	MG	A	1886	1/1	1.04	-	121,121,121,121	0
26	MG	A	1647	1/1	0.26	-	104,104,104,104	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1619	1/1	0.13	-	64,64,64,64	0
26	MG	A	1602	1/1	0.41	-	66,66,66,66	1
26	MG	A	1782	1/1	0.69	-	87,87,87,87	0
26	MG	A	1762	1/1	0.43	-	82,82,82,82	0
26	MG	A	1645	1/1	0.44	-	227,227,227,227	0
26	MG	D	304	1/1	0.43	-	455,455,455,455	0
26	MG	A	1620	1/1	0.27	-	166,166,166,166	0
27	SRY	A	1928	40/40	0.16	-	70,100,124,130	0
26	MG	A	1610	1/1	0.24	-	81,81,81,81	0
26	MG	D	303	1/1	0.20	-	104,104,104,104	0
26	MG	P	102	1/1	0.34	-	101,101,101,101	0
26	MG	A	1913	1/1	0.81	-	108,108,108,108	0
26	MG	A	1695	1/1	0.04	-	176,176,176,176	0
26	MG	A	1922	1/1	0.41	-	110,110,110,110	0
26	MG	A	1927	1/1	0.24	-	112,112,112,112	0
26	MG	A	1760	1/1	0.68	-	88,88,88,88	0
26	MG	A	1726	1/1	0.14	-	98,98,98,98	0
26	MG	A	1677	1/1	0.23	-	235,235,235,235	0
26	MG	A	1722	1/1	0.64	-	95,95,95,95	0
26	MG	A	1772	1/1	0.12	-	93,93,93,93	0
26	MG	A	1826	1/1	0.05	-	395,395,395,395	0
26	MG	A	1810	1/1	0.21	-	474,474,474,474	0
26	MG	A	1879	1/1	0.22	-	438,438,438,438	0
26	MG	A	1850	1/1	0.04	-	236,236,236,236	0
26	MG	A	1861	1/1	0.14	-	443,443,443,443	0
26	MG	A	1643	1/1	0.10	-	71,71,71,71	0
26	MG	A	1789	1/1	0.09	-	378,378,378,378	0
26	MG	A	1733	1/1	0.19	-	70,70,70,70	0
26	MG	A	1875	1/1	0.15	-	460,460,460,460	0
26	MG	A	1735	1/1	0.15	-	106,106,106,106	0
26	MG	A	1795	1/1	1.02	-	359,359,359,359	0
26	MG	A	1874	1/1	0.08	-	467,467,467,467	0
26	MG	A	1909	1/1	0.27	-	93,93,93,93	0
26	MG	A	1704	1/1	0.15	-	104,104,104,104	0
26	MG	A	1665	1/1	0.72	-	117,117,117,117	0
26	MG	A	1749	1/1	0.13	-	84,84,84,84	0
26	MG	A	1672	1/1	0.09	-	166,166,166,166	0
26	MG	A	1618	1/1	0.48	-	130,130,130,130	0
26	MG	A	1758	1/1	0.49	-	101,101,101,101	0
26	MG	A	1651	1/1	0.17	-	110,110,110,110	0
26	MG	A	1798	1/1	0.05	-	464,464,464,464	0
26	MG	A	1847	1/1	0.17	-	441,441,441,441	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1805	1/1	0.05	-	331,331,331,331	0
26	MG	A	1714	1/1	0.21	-	108,108,108,108	0
26	MG	A	1840	1/1	0.31	-	467,467,467,467	1
26	MG	A	1683	1/1	0.06	-	154,154,154,154	0
26	MG	A	1703	1/1	0.15	-	97,97,97,97	0
26	MG	A	1625	1/1	0.07	-	113,113,113,113	0
26	MG	A	1614	1/1	0.21	-	285,285,285,285	0
26	MG	A	1856	1/1	0.06	-	478,478,478,478	0
26	MG	A	1867	1/1	0.61	-	413,413,413,413	1
26	MG	A	1747	1/1	0.13	-	115,115,115,115	0
26	MG	S	102	1/1	0.16	-	106,106,106,106	0
26	MG	A	1707	1/1	0.40	-	135,135,135,135	0
26	MG	A	1718	1/1	0.20	-	326,326,326,326	0
26	MG	A	1631	1/1	0.09	-	158,158,158,158	0
26	MG	A	1713	1/1	0.15	-	211,211,211,211	0
26	MG	A	1877	1/1	0.27	-	456,456,456,456	1
26	MG	A	1684	1/1	0.14	-	109,109,109,109	0
26	MG	A	1924	1/1	0.16	-	131,131,131,131	0
26	MG	A	1635	1/1	0.13	-	74,74,74,74	0
26	MG	A	1697	1/1	0.19	-	135,135,135,135	0
26	MG	A	1744	1/1	0.32	-	82,82,82,82	0
26	MG	A	1792	1/1	0.13	-	415,415,415,415	0
26	MG	D	302	1/1	0.39	-	104,104,104,104	0
26	MG	A	1759	1/1	0.86	-	86,86,86,86	0
26	MG	A	1691	1/1	0.16	-	186,186,186,186	0
26	MG	A	1646	1/1	0.18	-	136,136,136,136	0
26	MG	A	1784	1/1	0.19	-	122,122,122,122	0
26	MG	A	1802	1/1	0.12	-	355,355,355,355	0
26	MG	A	1709	1/1	0.23	-	245,245,245,245	0
26	MG	A	1734	1/1	0.60	-	124,124,124,124	0
26	MG	A	1889	1/1	0.29	-	95,95,95,95	0
26	MG	A	1761	1/1	0.49	-	105,105,105,105	0
26	MG	A	1673	1/1	0.88	-	91,91,91,91	0
26	MG	A	1777	1/1	0.15	-	94,94,94,94	0
26	MG	A	1681	1/1	0.04	-	256,256,256,256	0
26	MG	A	1903	1/1	0.17	-	78,78,78,78	0
26	MG	A	1732	1/1	0.24	-	100,100,100,100	0
26	MG	A	1721	1/1	0.24	-	92,92,92,92	0
26	MG	A	1825	1/1	0.17	-	436,436,436,436	0
26	MG	A	1653	1/1	0.21	-	96,96,96,96	0
26	MG	A	1752	1/1	0.25	-	91,91,91,91	0
26	MG	A	1841	1/1	0.18	-	496,496,496,496	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
26	MG	A	1670	1/1	0.28	-	234,234,234,234	0
26	MG	A	1754	1/1	0.16	-	85,85,85,85	0
26	MG	A	1813	1/1	0.30	-	471,471,471,471	0
26	MG	A	1692	1/1	0.20	-	182,182,182,182	0
26	MG	A	1884	1/1	1.43	-	119,119,119,119	0
26	MG	A	1764	1/1	0.34	-	78,78,78,78	0
26	MG	A	1706	1/1	0.17	-	190,190,190,190	0
26	MG	A	1719	1/1	0.07	-	262,262,262,262	0
26	MG	A	1705	1/1	0.13	-	74,74,74,74	0
26	MG	A	1766	1/1	0.59	-	105,105,105,105	0
26	MG	A	1842	1/1	0.59	-	488,488,488,488	0
26	MG	A	1655	1/1	0.38	-	125,125,125,125	0
26	MG	A	1892	1/1	0.55	-	92,92,92,92	0
26	MG	A	1854	1/1	0.33	-	418,418,418,418	0
26	MG	A	1769	1/1	0.62	-	127,127,127,127	0
26	MG	A	1891	1/1	0.16	-	127,127,127,127	0
26	MG	A	1833	1/1	0.12	-	335,335,335,335	0
26	MG	A	1652	1/1	0.09	-	72,72,72,72	0
26	MG	A	1687	1/1	0.14	-	106,106,106,106	0
26	MG	A	1831	1/1	0.22	-	395,395,395,395	0
26	MG	A	1799	1/1	0.23	-	426,426,426,426	0
26	MG	A	1790	1/1	0.15	-	444,444,444,444	0
26	MG	A	1629	1/1	0.39	-	223,223,223,223	0

6.5 Other polymers ⓘ

There are no such residues in this entry.