



# wwPDB X-ray Structure Validation Summary Report

Feb 27, 2014 – 02:10 PM GMT

PDB ID : 1KD1  
Title : Co-crystal Structure of Spiramycin bound to the 50S Ribosomal Subunit of *Haloarcula marismortui*  
Authors : Hansen, J.L.; Ippolito, J.A.; Ban, N.; Nissen, P.; Moore, P.B.; Steitz, T.A.  
Deposited on : 2001-11-12  
Resolution : 3.00 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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

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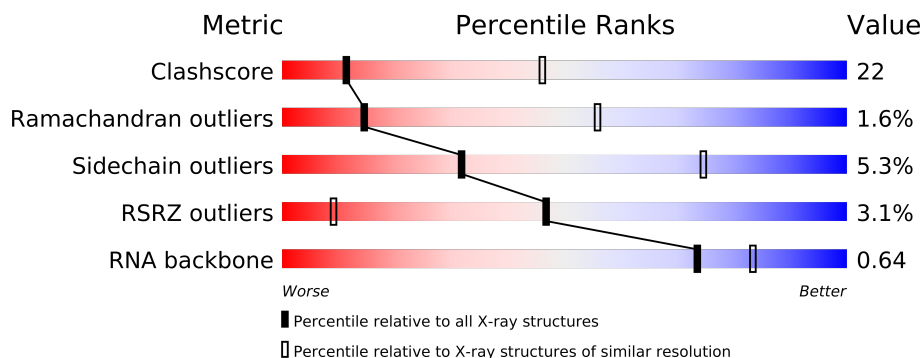
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.00 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))
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)
RNA backbone	1838	1070 (3.50-2.50)

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  $\geq 3$ , 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	2922	
2	B	122	
3	C	239	
4	D	337	
5	E	246	
6	F	176	
7	G	177	
8	H	119	
9	I	348	
10	J	167	
11	K	145	
12	L	132	
13	M	164	
14	N	194	

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Mol	Chain	Length	Quality of chain
15	O	186	
16	P	115	
17	Q	148	
18	R	95	
19	S	154	
20	T	84	
21	U	119	
22	V	66	
23	W	70	
24	X	154	
25	Y	91	
26	Z	240	
27	1	73	
28	2	56	
29	3	48	
30	4	92	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
31	SPR	A	9001	-	X
32	MG	A	8024	-	X
32	MG	A	8049	-	X
32	MG	A	8070	-	X
32	MG	A	8082	-	X
32	MG	A	8090	-	X
32	MG	A	8092	-	X
32	MG	A	8093	-	X
32	MG	A	8094	-	X
32	MG	A	8097	-	X
32	MG	A	8101	-	X
32	MG	A	8102	-	X
32	MG	A	8103	-	X
32	MG	A	8104	-	X
32	MG	A	8114	-	X
32	MG	A	8118	-	X
32	MG	A	8119	-	X
32	MG	Z	8109	-	X
33	NA	A	8306	-	X
33	NA	A	8308	-	X
33	NA	A	8313	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
33	NA	A	8316	-	X
33	NA	A	8318	-	X
33	NA	A	8321	-	X
33	NA	A	8322	-	X
33	NA	A	8323	-	X
33	NA	A	8328	-	X
33	NA	A	8329	-	X
33	NA	A	8331	-	X
33	NA	A	8332	-	X
33	NA	A	8336	-	X
33	NA	A	8340	-	X
33	NA	A	8342	-	X
33	NA	A	8350	-	X
33	NA	A	8352	-	X
33	NA	A	8355	-	X
33	NA	A	8356	-	X
33	NA	A	8359	-	X
33	NA	A	8360	-	X
33	NA	A	8362	-	X
33	NA	A	8363	-	X
33	NA	A	8364	-	X
33	NA	A	8365	-	X
33	NA	A	8366	-	X
33	NA	A	8367	-	X
33	NA	A	8370	-	X
33	NA	A	8371	-	X
33	NA	A	8372	-	X
33	NA	A	8373	-	X
33	NA	A	8374	-	X
33	NA	A	8376	-	X
33	NA	A	8377	-	X
33	NA	A	8378	-	X
33	NA	A	8379	-	X
33	NA	A	8382	-	X
33	NA	A	8384	-	X
33	NA	A	8385	-	X
33	NA	B	8383	-	X
33	NA	S	8337	-	X
33	NA	S	8386	-	X
33	NA	T	8312	-	X
34	CL	4	8504	-	X
34	CL	A	8503	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
34	CL	A	8505	-	X
34	CL	A	8515	-	X
34	CL	A	8517	-	X
34	CL	A	8522	-	X
34	CL	C	8509	-	X
34	CL	D	8519	-	X
34	CL	R	8511	-	X
35	K	A	8603	-	X
36	CD	4	8404	-	X
36	CD	P	8405	-	X

## 2 Entry composition

There are 37 unique types of molecules in this entry. The entry contains 98587 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 23S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	2754	Total	C	N	O	P	0	0	0
			59017	26346	10878	19048	2745			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	560	C	U	CONFLICT	? 3377779

- Molecule 2 is a RNA chain called 5S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	122	Total	C	N	O	P	0	0	0
			2600	1160	472	847	121			

- Molecule 3 is a protein called RIBOSOMAL PROTEIN L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	237	Total	C	N	O	S	0	0	0
			1754	1072	352	325	5			

- Molecule 4 is a protein called RIBOSOMAL PROTEIN L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	337	Total	C	N	O	S	0	0	0
			2624	1616	493	510	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	PRO	DELETION	UNP P20279
D	310	ARG	PHE	CONFLICT	UNP P20279

- Molecule 5 is a protein called RIBOSOMAL PROTEIN L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	246	Total	C	N	O	S	0	0	0
			1858	1131	344	382	1			

- Molecule 6 is a protein called RIBOSOMAL PROTEIN L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	140	Total	C	N	O	S	0	0	0
			1094	685	195	210	4			

- Molecule 7 is a protein called RIBOSOMAL PROTEIN L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	172	Total	C	N	O	S	0	0	0
			1357	840	224	289	4			

- Molecule 8 is a protein called RIBOSOMAL PROTEIN L7AE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	119	Total	C	N	O	S	0	0	0
			885	552	141	191	1			

- Molecule 9 is a protein called RIBOSOMAL PROTEIN L10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	29	Total	C	N	O	S	0	0	0
			240	149	39	51	1			

- Molecule 10 is a protein called RIBOSOMAL PROTEIN L10E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	156	Total	C	N	O	S	0	0	0
			1215	766	233	212	4			

- Molecule 11 is a protein called RIBOSOMAL PROTEIN L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	142	Total	C	N	O	S	0	0	0
			1119	696	199	221	3			

- Molecule 12 is a protein called RIBOSOMAL PROTEIN L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	132	Total	C	N	O	S	0	0	0
			993	609	189	191	4			

- Molecule 13 is a protein called RIBOSOMAL PROTEIN L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	145	Total	C	N	O		0	0	0
			1114	668	222	224				

- Molecule 14 is a protein called RIBOSOMAL PROTEIN L15E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	194	Total	C	N	O	S	0	0	0
			1605	988	346	266	5			

- Molecule 15 is a protein called RIBOSOMAL PROTEIN L18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	O	186	Total	C	N	O	S	0	0	0
			1444	895	262	285	2			

- Molecule 16 is a protein called RIBOSOMAL PROTEIN L18E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	P	115	Total	C	N	O		0	0	0
			864	529	161	174				

- Molecule 17 is a protein called RIBOSOMAL PROTEIN L19E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	Q	143	Total	C	N	O		0	0	0
			1133	680	230	223				

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	71	LYS	TYR	CONFLICT	UNP P14119

- Molecule 18 is a protein called RIBOSOMAL PROTEIN L21E.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	R	95	Total	C	N	O			
			734	450	141	143	0	0	0

- Molecule 19 is a protein called RIBOSOMAL PROTEIN L22.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
19	S	150	Total	C	N	O	S		
			1149	713	209	223	4	0	0

- Molecule 20 is a protein called RIBOSOMAL PROTEIN L23.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
20	T	81	Total	C	N	O	S		
			641	389	111	138	3	0	0

- Molecule 21 is a protein called RIBOSOMAL PROTEIN L24.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	U	119	Total	C	N	O			
			949	568	180	201		0	0

- Molecule 22 is a protein called RIBOSOMAL PROTEIN L24E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
22	V	53	Total	C	N	O	S		
			410	244	75	86	5	0	0

- Molecule 23 is a protein called RIBOSOMAL PROTEIN L29.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
23	W	65	Total	C	N	O	S		
			499	304	94	100	1	0	0

- Molecule 24 is a protein called RIBOSOMAL PROTEIN L30.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
24	X	154	Total	C	N	O	S		
			1195	737	209	243	6	0	0

- Molecule 25 is a protein called RIBOSOMAL PROTEIN L31E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	Y	82	Total	C	N	O	S	0	0	0
			654	402	129	122	1			

- Molecule 26 is a protein called RIBOSOMAL PROTEIN L32E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	Z	142	Total	C	N	O		0	0	0
			1130	686	228	216				

- Molecule 27 is a protein called RIBOSOMAL PROTEIN L37Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	1	73	Total	C	N	O	S	0	0	0
			563	359	111	86	7			

- Molecule 28 is a protein called RIBOSOMAL PROTEIN L37E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	2	56	Total	C	N	O	S	0	0	0
			430	258	86	82	4			

- Molecule 29 is a protein called RIBOSOMAL PROTEIN L39E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	3	46	Total	C	N	O	S	0	0	0
			393	238	86	68	1			

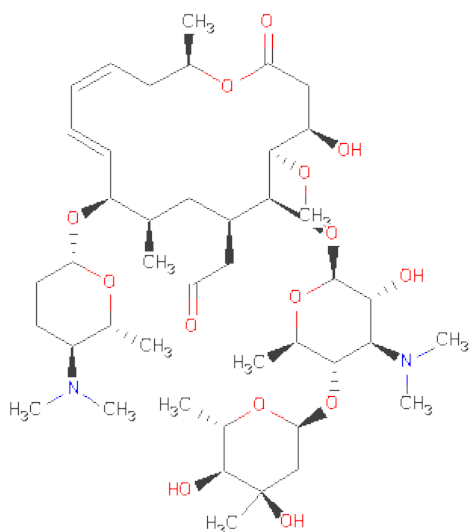
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	?	-	ARG	DELETION	UNP P22452

- Molecule 30 is a protein called RIBOSOMAL PROTEIN L44E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	4	92	Total	C	N	O	S	0	0	0
			755	458	153	137	7			

- Molecule 31 is SPIRAMYCIN I (three-letter code: SPR) (formula: C<sub>43</sub>H<sub>74</sub>N<sub>2</sub>O<sub>14</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
31	A	1	Total	C	N	O	0	0
			59	43	2	14		

- Molecule 32 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
32	1	1	Total	Mg	0	0
			1	1		
32	B	1	Total	Mg	0	0
			1	1		
32	C	1	Total	Mg	0	0
			1	1		
32	Z	1	Total	Mg	0	0
			1	1		
32	A	112	Total	Mg	0	0
			112	112		
32	4	1	Total	Mg	0	0
			1	1		
32	U	1	Total	Mg	0	0
			1	1		
32	L	1	Total	Mg	0	0
			1	1		

- Molecule 33 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
33	J	1	Total Na 1 1	0	0
33	K	1	Total Na 1 1	0	0
33	E	1	Total Na 1 1	0	0
33	B	2	Total Na 2 2	0	0
33	C	1	Total Na 1 1	0	0
33	A	73	Total Na 73 73	0	0
33	T	1	Total Na 1 1	0	0
33	N	1	Total Na 1 1	0	0
33	R	1	Total Na 1 1	0	0
33	S	2	Total Na 2 2	0	0
33	M	1	Total Na 1 1	0	0

- Molecule 34 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
34	P	1	Total Cl 1 1	0	0
34	D	1	Total Cl 1 1	0	0
34	K	3	Total Cl 3 3	0	0
34	C	1	Total Cl 1 1	0	0
34	Z	1	Total Cl 1 1	0	0
34	A	9	Total Cl 9 9	0	0
34	4	1	Total Cl 1 1	0	0
34	N	1	Total Cl 1 1	0	0
34	O	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
34	R	1	Total 1	Cl 1	0	0
34	S	1	Total 1	Cl 1	0	0
34	M	1	Total 1	Cl 1	0	0

- Molecule 35 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
35	A	3	Total 3	K 3	0	0

- Molecule 36 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
36	P	1	Total 1	Cd 1	0	0
36	2	1	Total 1	Cd 1	0	0
36	1	1	Total 1	Cd 1	0	0
36	4	1	Total 1	Cd 1	0	0
36	V	1	Total 1	Cd 1	0	0

- Molecule 37 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	A	5910	Total 5910	O 5910	0	0
37	B	142	Total 142	O 142	0	0
37	C	126	Total 126	O 126	0	0
37	D	150	Total 150	O 150	0	0
37	E	169	Total 169	O 169	0	0
37	F	51	Total 51	O 51	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	G	42	Total 42	O 42	0	0
37	H	26	Total 26	O 26	0	0
37	I	21	Total 21	O 21	0	0
37	J	78	Total 78	O 78	0	0
37	K	54	Total 54	O 54	0	0
37	L	65	Total 65	O 65	0	0
37	M	79	Total 79	O 79	0	0
37	N	132	Total 132	O 132	0	0
37	O	69	Total 69	O 69	0	0
37	P	45	Total 45	O 45	0	0
37	Q	65	Total 65	O 65	0	0
37	R	55	Total 55	O 55	0	0
37	S	83	Total 83	O 83	0	0
37	T	35	Total 35	O 35	0	0
37	U	39	Total 39	O 39	0	0
37	V	25	Total 25	O 25	0	0
37	W	15	Total 15	O 15	0	0
37	X	70	Total 70	O 70	0	0
37	Y	25	Total 25	O 25	0	0
37	Z	94	Total 94	O 94	0	0
37	1	41	Total 41	O 41	0	0

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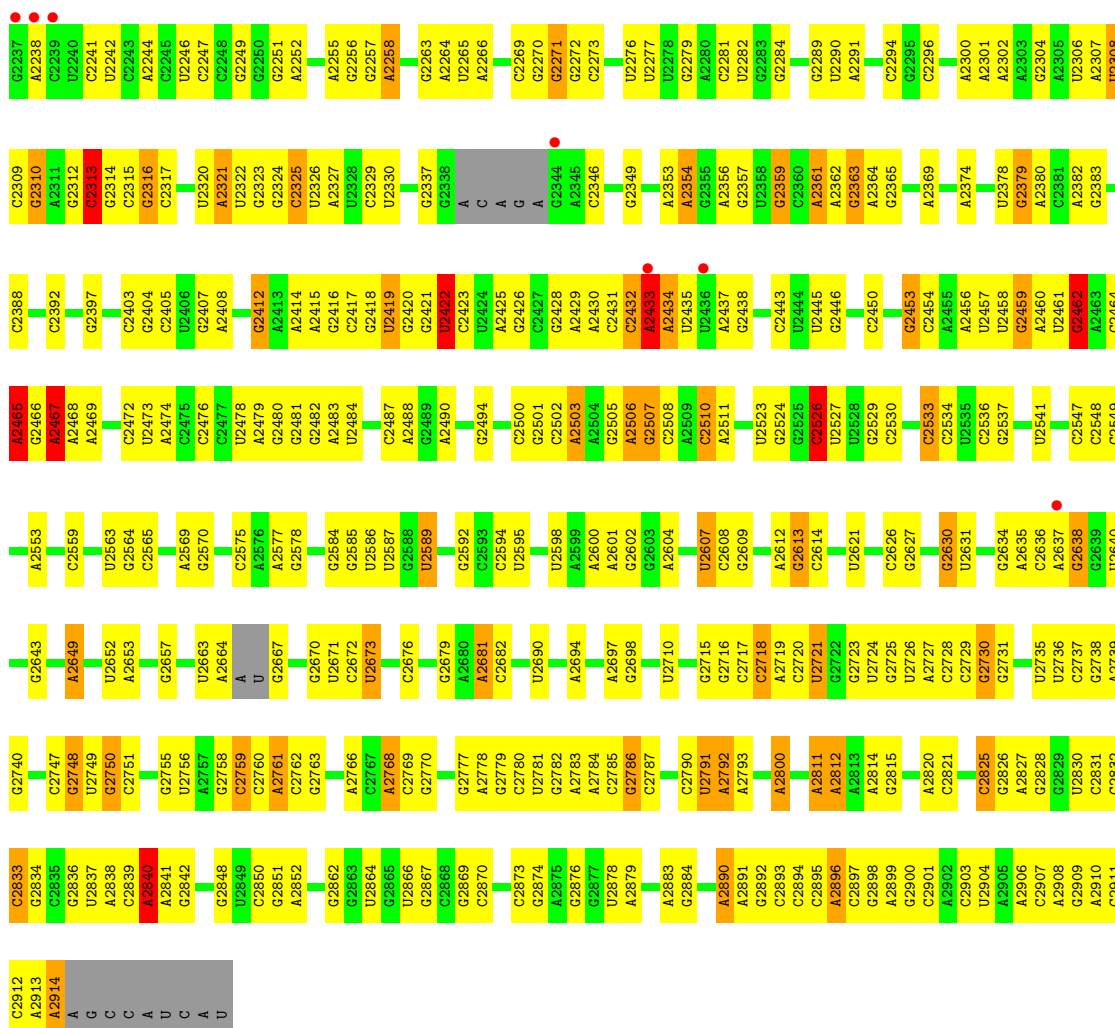
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	2	55	Total 55	O 55	0	0
37	3	42	Total 42	O 42	0	0
37	4	73	Total 73	O 73	0	0



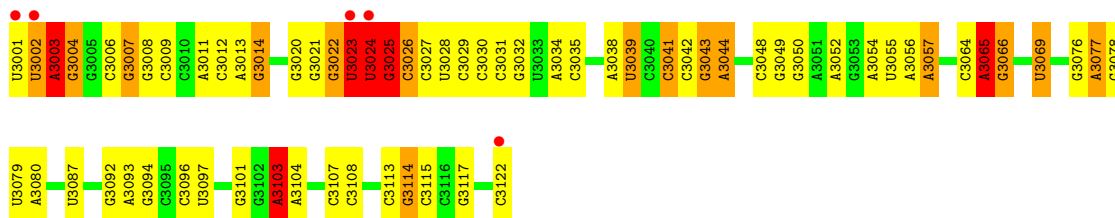






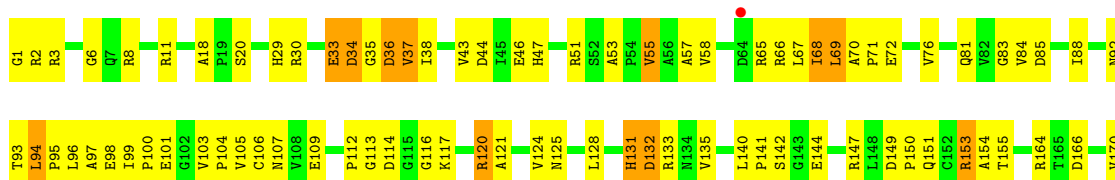
• Molecule 2: 5S RRNA

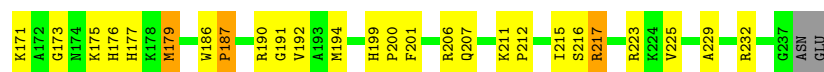
Chain B:



• Molecule 3: RIBOSOMAL PROTEIN L2

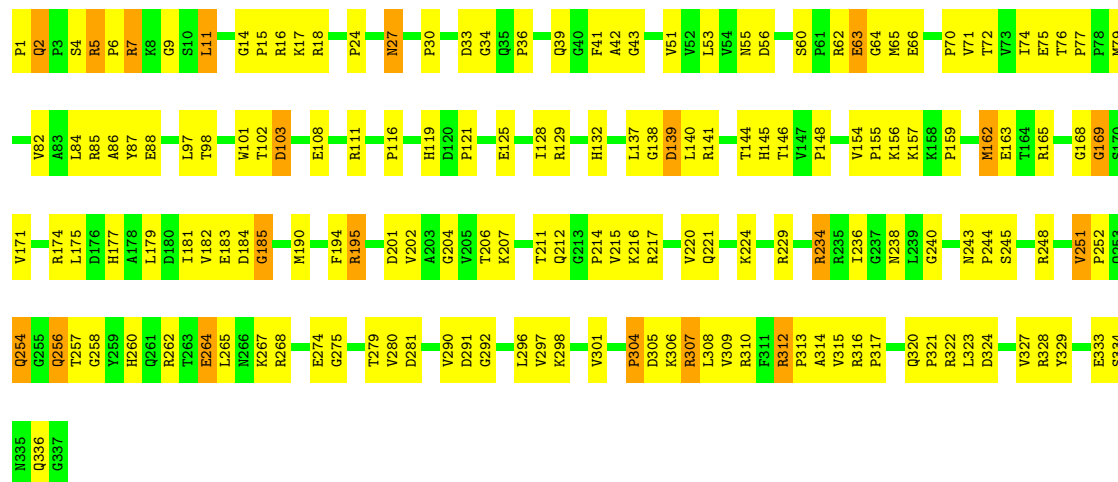
Chain C:





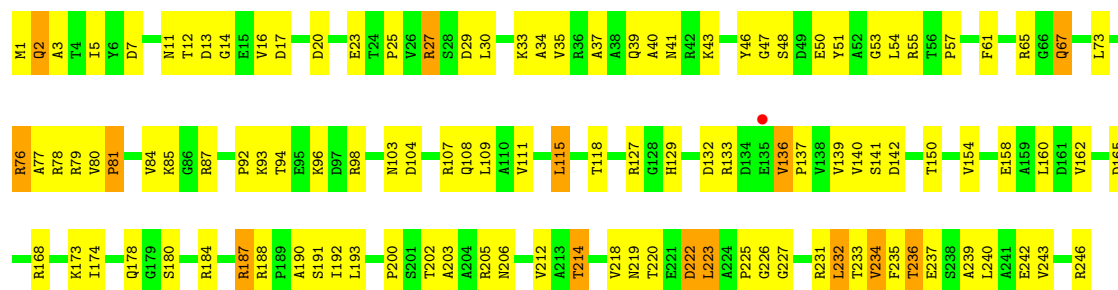
### • Molecule 4: RIBOSOMAL PROTEIN L3

Chain D:



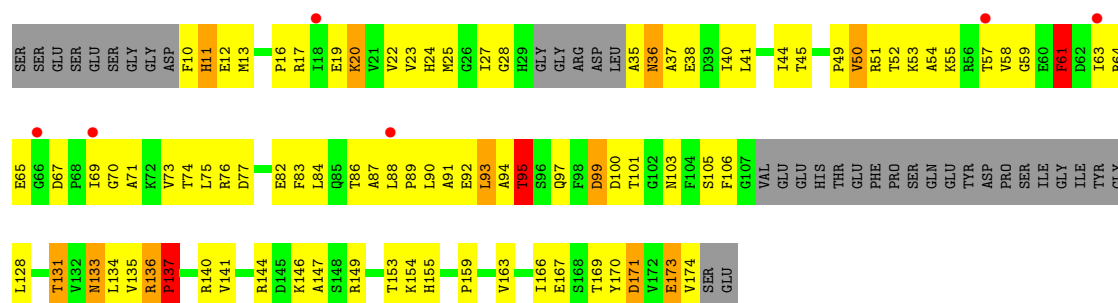
### • Molecule 5: RIBOSOMAL PROTEIN L4

Chain E:



### • Molecule 6: RIBOSOMAL PROTEIN L5

Chain F:



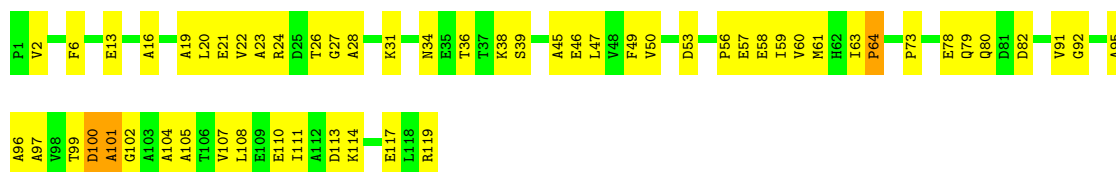
### • Molecule 7: RIBOSOMAL PROTEIN L6

Chain G:



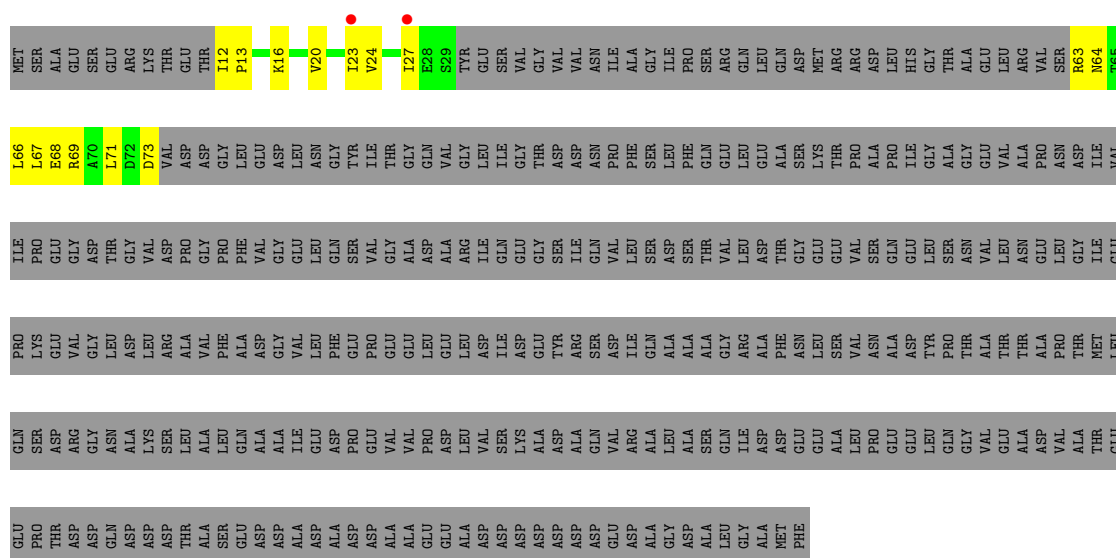
• Molecule 8: RIBOSOMAL PROTEIN L7AE

Chain H:



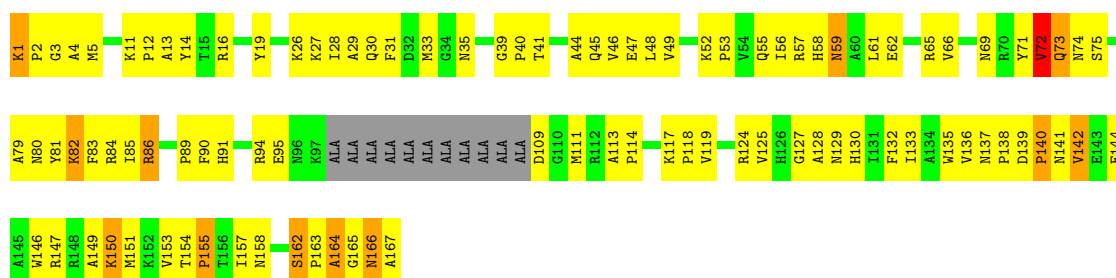
• Molecule 9: RIBOSOMAL PROTEIN L10

Chain I:



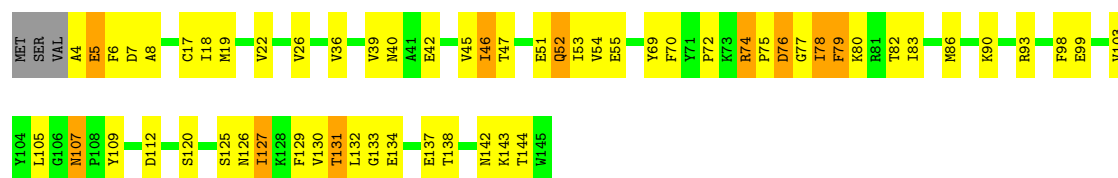
• Molecule 10: RIBOSOMAL PROTEIN L10E

Chain J:



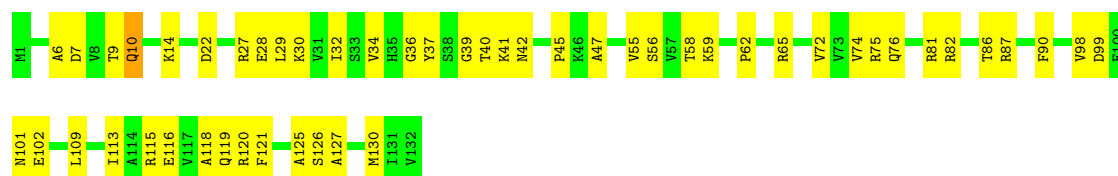
• Molecule 11: RIBOSOMAL PROTEIN L13

Chain K: 



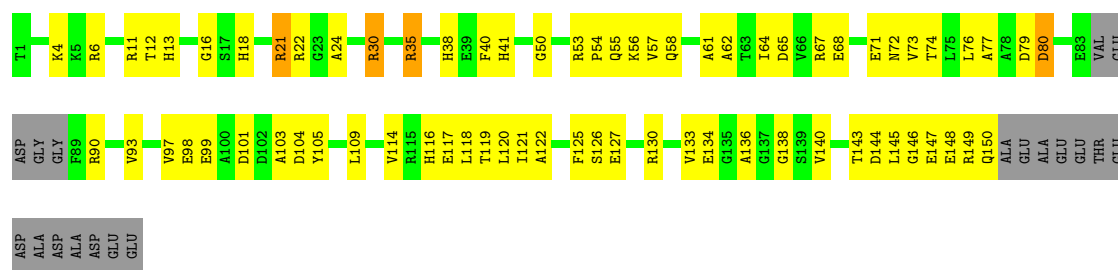
• Molecule 12: RIBOSOMAL PROTEIN L14

Chain L: 



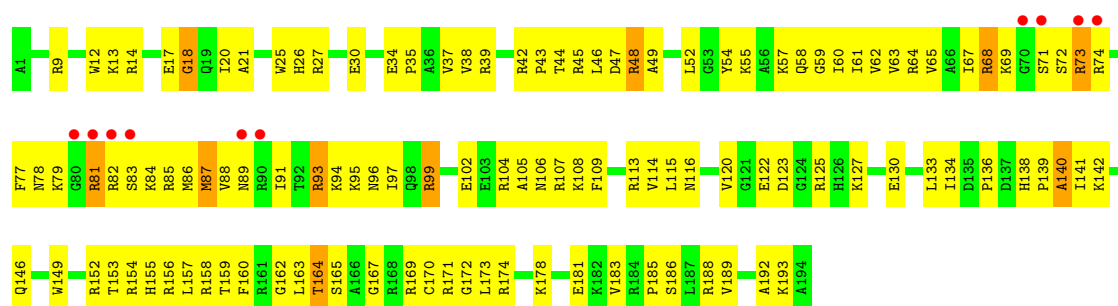
• Molecule 13: RIBOSOMAL PROTEIN L15

Chain M: 



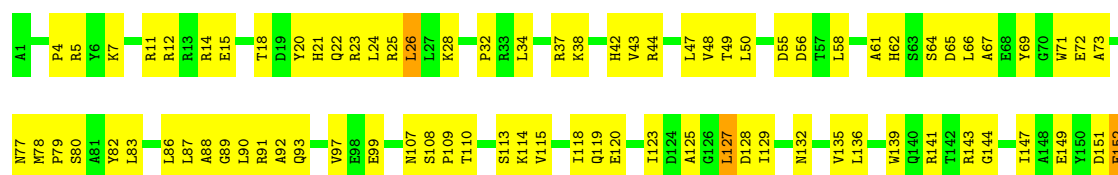
• Molecule 14: RIBOSOMAL PROTEIN L15E

Chain N: 



• Molecule 15: RIBOSOMAL PROTEIN L18

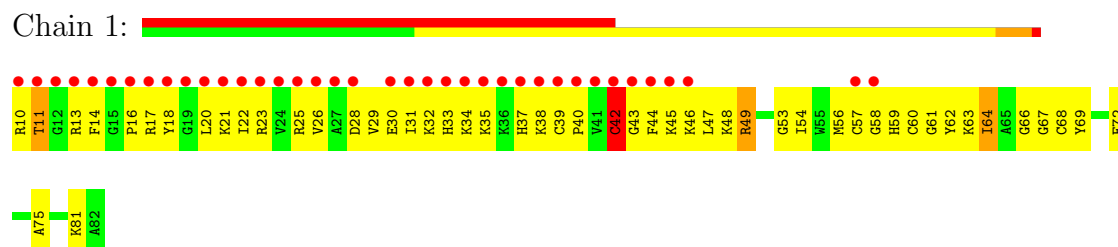
Chain O: 



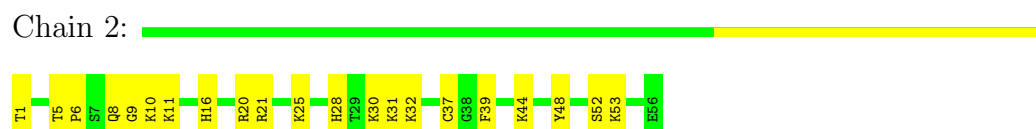




- Molecule 27: RIBOSOMAL PROTEIN L37Ae



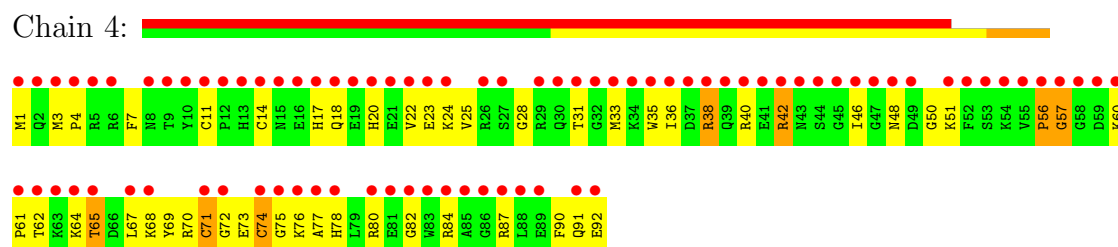
- Molecule 28: RIBOSOMAL PROTEIN L37E



- Molecule 29: RIBOSOMAL PROTEIN L39E



- Molecule 30: RIBOSOMAL PROTEIN L44E





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	212.90Å 300.47Å 575.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 3.00 50.07 – 2.99	Depositor EDS
% Data completeness (in resolution range)	91.7 (19.99-3.00) 91.1 (50.07-2.99)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	0.15	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.46 (at 3.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.220 , 0.269 0.220 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	50.9	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 30.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	1 of 363802 reflections (0.000%)	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	98587	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.66% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: MG, CL, NA, SPR, CD, K

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.62	12/66076 (0.0%)	0.79	45/103052 (0.0%)
2	B	0.90	12/2905 (0.4%)	0.98	17/4528 (0.4%)
3	C	0.53	0/1787	0.79	0/2409
4	D	0.52	0/2689	0.75	0/3652
5	E	0.54	0/1883	0.78	0/2551
6	F	0.43	0/1111	0.65	0/1498
7	G	0.47	0/1382	0.66	0/1880
8	H	0.44	0/896	0.64	0/1219
9	I	0.41	0/241	0.58	0/324
10	J	0.53	0/1246	0.83	1/1686 (0.1%)
11	K	0.52	0/1135	0.70	0/1530
12	L	0.51	0/1003	0.80	0/1351
13	M	0.49	0/1126	0.74	0/1504
14	N	0.67	0/1633	0.86	1/2180 (0.0%)
15	O	0.48	0/1473	0.76	0/1999
16	P	0.53	0/873	0.76	0/1181
17	Q	0.52	0/1143	0.67	0/1521
18	R	0.52	0/748	0.80	1/1005 (0.1%)
19	S	0.66	1/1172 (0.1%)	0.84	2/1578 (0.1%)
20	T	0.45	0/648	0.69	0/875
21	U	0.47	0/957	0.73	1/1289 (0.1%)
22	V	0.77	0/417	0.86	2/562 (0.4%)
23	W	0.42	0/502	0.63	0/675
24	X	0.54	0/1218	0.76	0/1655
25	Y	0.50	0/664	0.72	0/895
26	Z	0.53	0/1146	0.73	0/1536
27	1	0.85	0/575	0.87	1/763 (0.1%)
28	2	0.56	0/437	0.84	0/578
29	3	0.47	0/398	0.64	0/527
30	4	1.04	0/771	0.83	1/1024 (0.1%)
All	All	0.62	25/98255 (0.0%)	0.79	72/147027 (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
1	A	1	198
2	B	0	6
28	2	0	1
All	All	1	205

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2103	A	C5-C6	13.31	1.53	1.41
2	B	3025	G	O3'-P	11.56	1.75	1.61
2	B	3026	C	P-OP2	-10.89	1.30	1.49
2	B	3026	C	P-O5'	-9.81	1.50	1.59
2	B	3023	U	C2'-O2'	8.99	1.53	1.41

The worst 5 of 72 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1164	U	OP2-P-O3'	-18.78	63.89	105.20
1	A	1164	U	OP1-P-O3'	-18.34	64.85	105.20
1	A	2104	C	O5'-P-OP1	-14.12	92.99	105.70
2	B	3024	U	O5'-P-OP2	11.53	124.53	110.70
2	B	3026	C	O5'-P-OP2	-11.17	95.65	105.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	1563	G	C3'

5 of 205 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	G	Sidechain
1	A	138	U	Sidechain
1	A	146	U	Sidechain
1	A	75	U	Sidechain
1	A	99	A	Sidechain

## 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	59017	0	29802	1290	0
2	B	2600	0	1326	80	0
3	C	1754	0	1763	132	0
4	D	2624	0	2533	190	0
5	E	1858	0	1816	149	0
6	F	1094	0	1085	135	0
7	G	1357	0	1266	85	0
8	H	885	0	854	63	0
9	I	240	0	231	21	0
10	J	1215	0	1215	175	0
11	K	1119	0	1098	70	0
12	L	993	0	1027	67	0
13	M	1114	0	1072	72	0
14	N	1605	0	1676	194	0
15	O	1444	0	1401	143	0
16	P	864	0	873	37	0
17	Q	1133	0	1127	53	0
18	R	734	0	727	30	0
19	S	1149	0	1122	60	0
20	T	641	0	605	23	0
21	U	949	0	923	59	0
22	V	410	0	368	45	0
23	W	499	0	511	33	0
24	X	1195	0	1137	99	0
25	Y	654	0	653	51	0
26	Z	1130	0	1133	71	0
27	1	563	0	601	80	0
28	2	430	0	426	27	0
29	3	393	0	406	27	0
30	4	755	0	732	58	0
31	A	59	0	73	9	0
32	1	1	0	0	0	0
32	4	1	0	0	0	0
32	A	112	0	0	5	0
32	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	C	1	0	0	0	0
32	L	1	0	0	0	0
32	U	1	0	0	0	0
32	Z	1	0	0	0	0
33	A	73	0	0	1	0
33	B	2	0	0	0	0
33	C	1	0	0	0	0
33	E	1	0	0	0	0
33	J	1	0	0	0	0
33	K	1	0	0	0	0
33	M	1	0	0	0	0
33	N	1	0	0	0	0
33	R	1	0	0	0	0
33	S	2	0	0	0	0
33	T	1	0	0	0	0
34	4	1	0	0	0	0
34	A	9	0	0	2	0
34	C	1	0	0	0	0
34	D	1	0	0	0	0
34	K	3	0	0	2	0
34	M	1	0	0	0	0
34	N	1	0	0	1	0
34	O	1	0	0	3	0
34	P	1	0	0	0	0
34	R	1	0	0	0	0
34	S	1	0	0	0	0
34	Z	1	0	0	0	0
35	A	3	0	0	0	0
36	1	1	0	0	0	0
36	2	1	0	0	0	0
36	4	1	0	0	0	0
36	P	1	0	0	0	0
36	V	1	0	0	0	0
37	1	41	0	0	13	0
37	2	55	0	0	5	0
37	3	42	0	0	5	0
37	4	73	0	0	7	0
37	A	5910	0	0	300	0
37	B	142	0	0	16	0
37	C	126	0	0	23	0
37	D	150	0	0	28	0
37	E	169	0	0	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
37	F	51	0	0	22	0
37	G	42	0	0	13	0
37	H	26	0	0	9	0
37	I	21	0	0	5	0
37	J	78	0	0	26	0
37	K	54	0	0	8	0
37	L	65	0	0	12	0
37	M	79	0	0	18	0
37	N	132	0	0	36	0
37	O	69	0	0	23	0
37	P	45	0	0	9	0
37	Q	65	0	0	4	0
37	R	55	0	0	6	0
37	S	83	0	0	11	0
37	T	35	0	0	3	0
37	U	39	0	0	4	0
37	V	25	0	0	8	0
37	W	15	0	0	2	0
37	X	70	0	0	10	0
37	Y	25	0	0	11	0
37	Z	94	0	0	18	0
All	All	98587	0	59582	3325	0

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 22.

The worst 5 of 3325 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
30:4:33:MET:SD	30:4:33:MET:CE	2.03	1.47
5:E:236:THR:HG22	5:E:239:ALA:H	1.09	1.15
1:A:2121:G:OP2	37:A:3494:HOH:O	1.64	1.15
1:A:2122:C:OP2	37:A:6549:HOH:O	1.64	1.15
1:A:1134:G:H4'	10:J:151:MET:HE1	1.28	1.12

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
3	C	235/239 (98%)	205 (87%)	26 (11%)	4 (2%)	14	54
4	D	335/337 (99%)	307 (92%)	21 (6%)	7 (2%)	11	47
5	E	244/246 (99%)	225 (92%)	18 (7%)	1 (0%)	43	87
6	F	134/176 (76%)	95 (71%)	27 (20%)	12 (9%)	1	5
7	G	170/177 (96%)	158 (93%)	12 (7%)	0	100	100
8	H	117/119 (98%)	103 (88%)	12 (10%)	2 (2%)	14	54
9	I	25/348 (7%)	24 (96%)	1 (4%)	0	100	100
10	J	152/167 (91%)	129 (85%)	17 (11%)	6 (4%)	5	26
11	K	140/145 (97%)	131 (94%)	5 (4%)	4 (3%)	7	35
12	L	130/132 (98%)	117 (90%)	11 (8%)	2 (2%)	15	58
13	M	141/164 (86%)	118 (84%)	20 (14%)	3 (2%)	11	47
14	N	192/194 (99%)	172 (90%)	18 (9%)	2 (1%)	22	70
15	O	184/186 (99%)	165 (90%)	13 (7%)	6 (3%)	6	32
16	P	113/115 (98%)	110 (97%)	3 (3%)	0	100	100
17	Q	141/148 (95%)	138 (98%)	2 (1%)	1 (1%)	30	78
18	R	93/95 (98%)	87 (94%)	5 (5%)	1 (1%)	21	67
19	S	148/154 (96%)	138 (93%)	10 (7%)	0	100	100
20	T	79/84 (94%)	72 (91%)	7 (9%)	0	100	100
21	U	117/119 (98%)	108 (92%)	9 (8%)	0	100	100
22	V	51/66 (77%)	47 (92%)	4 (8%)	0	100	100
23	W	63/70 (90%)	57 (90%)	4 (6%)	2 (3%)	6	33
24	X	152/154 (99%)	145 (95%)	6 (4%)	1 (1%)	30	78
25	Y	80/91 (88%)	71 (89%)	7 (9%)	2 (2%)	9	40
26	Z	140/240 (58%)	138 (99%)	2 (1%)	0	100	100
27	1	71/73 (97%)	63 (89%)	7 (10%)	1 (1%)	16	60
28	2	54/56 (96%)	50 (93%)	4 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
29	3	42/48 (88%)	41 (98%)	1 (2%)	0	100	100
30	4	90/92 (98%)	85 (94%)	3 (3%)	2 (2%)	10	45
All	All	3633/4235 (86%)	3299 (91%)	275 (8%)	59 (2%)	14	56

5 of 59 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	139	ASP
6	F	93	LEU
6	F	95	THR
6	F	137	PRO
6	F	173	GLU

### 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	
3	C	179/181 (99%)	166 (93%)	13 (7%)	20	59
4	D	282/282 (100%)	264 (94%)	18 (6%)	25	66
5	E	193/193 (100%)	178 (92%)	15 (8%)	18	55
6	F	117/147 (80%)	106 (91%)	11 (9%)	13	44
7	G	152/155 (98%)	147 (97%)	5 (3%)	50	88
8	H	92/92 (100%)	91 (99%)	1 (1%)	84	97
9	I	27/283 (10%)	27 (100%)	0	100	100
10	J	122/122 (100%)	111 (91%)	11 (9%)	14	47
11	K	118/121 (98%)	107 (91%)	11 (9%)	13	45
12	L	106/106 (100%)	103 (97%)	3 (3%)	56	91
13	M	112/126 (89%)	108 (96%)	4 (4%)	47	86
14	N	166/166 (100%)	157 (95%)	9 (5%)	31	74
15	O	149/149 (100%)	144 (97%)	5 (3%)	49	88
16	P	93/93 (100%)	90 (97%)	3 (3%)	51	89
17	Q	113/116 (97%)	110 (97%)	3 (3%)	57	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
18	R	79/79 (100%)	75 (95%)	4 (5%)	33	76
19	S	117/121 (97%)	112 (96%)	5 (4%)	40	82
20	T	71/73 (97%)	71 (100%)	0	100	100
21	U	105/105 (100%)	103 (98%)	2 (2%)	69	94
22	V	44/52 (85%)	42 (96%)	2 (4%)	38	81
23	W	51/56 (91%)	50 (98%)	1 (2%)	68	94
24	X	130/130 (100%)	122 (94%)	8 (6%)	26	67
25	Y	66/73 (90%)	61 (92%)	5 (8%)	19	57
26	Z	120/195 (62%)	112 (93%)	8 (7%)	23	64
27	1	56/56 (100%)	50 (89%)	6 (11%)	10	35
28	2	46/46 (100%)	46 (100%)	0	100	100
29	3	42/44 (96%)	41 (98%)	1 (2%)	61	92
30	4	79/79 (100%)	73 (92%)	6 (8%)	19	57
All	All	3027/3441 (88%)	2867 (95%)	160 (5%)	32	74

5 of 160 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
10	J	155	PRO
13	M	35	ARG
27	1	11	THR
11	K	46	ILE
11	K	120	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 84 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	116	HIS
17	Q	73	HIS
28	2	16	HIS
14	N	58	GLN
14	N	176	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	2747/2922 (94%)	248 (9%)	40 (1%)
2	B	121/122 (99%)	15 (12%)	5 (4%)
All	All	2868/3044 (94%)	263 (9%)	45 (1%)

5 of 263 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	11	A
1	A	31	C
1	A	60	A
1	A	67	A
1	A	69	A

5 of 45 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	1377	C
1	A	1692	C
2	B	3023	U
1	A	1506	U
1	A	1856	C

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 235 ligands modelled in this entry, 234 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
31	SPR	A	9001	1	62,62,62	3.11	27 (43%)	89,89,89	2.89	34 (38%)

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
31	SPR	A	9001	1	-	0/61/113/113	0/3/4/4

The worst 5 of 27 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
31	A	9001	SPR	C14-C15	8.58	1.60	1.52
31	A	9001	SPR	O1C-C9	7.32	1.52	1.44
31	A	9001	SPR	C5C-C4C	6.98	1.64	1.53
31	A	9001	SPR	C3C-C4C	6.75	1.68	1.52
31	A	9001	SPR	C4C-N4C	6.11	1.63	1.48

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	A	9001	SPR	C8A-N3A-C7A	-8.69	82.41	110.40
31	A	9001	SPR	C7C-N4C-C4C	7.70	135.04	113.11
31	A	9001	SPR	C16-C15-C14	7.31	123.58	113.06
31	A	9001	SPR	C5A-C4A-C3A	-7.21	93.97	110.63
31	A	9001	SPR	O1C-C9-C8	6.65	124.98	108.33

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	2754/2922 (94%)	-0.23	34 (1%) 75 20	23, 51, 96, 143	0
2	B	122/122 (100%)	0.09	5 (4%) 35 7	37, 70, 100, 150	0
3	C	237/239 (99%)	-0.06	1 (0%) 90 41	32, 63, 96, 110	0
4	D	337/337 (100%)	-0.16	0 100 100	28, 57, 84, 95	0
5	E	246/246 (100%)	-0.19	1 (0%) 90 41	24, 50, 74, 85	0
6	F	140/176 (79%)	0.67	6 (4%) 34 7	60, 103, 122, 127	0
7	G	172/177 (97%)	0.11	0 100 100	43, 68, 92, 98	0
8	H	119/119 (100%)	0.15	0 100 100	59, 79, 102, 107	0
9	I	29/348 (8%)	0.81	2 (6%) 17 4	76, 94, 102, 104	0
10	J	156/167 (93%)	-0.04	0 100 100	35, 58, 85, 93	0
11	K	142/145 (97%)	-0.16	0 100 100	36, 50, 76, 84	0
12	L	132/132 (100%)	-0.12	0 100 100	35, 56, 78, 82	0
13	M	145/164 (88%)	0.20	0 100 100	31, 74, 108, 117	0
14	N	194/194 (100%)	0.05	10 (5%) 26 6	37, 55, 91, 98	0
15	O	186/186 (100%)	0.29	3 (1%) 68 16	48, 74, 112, 122	0
16	P	115/115 (100%)	-0.10	0 100 100	39, 59, 75, 79	0
17	Q	143/148 (96%)	0.02	0 100 100	38, 60, 76, 84	0
18	R	95/95 (100%)	-0.10	0 100 100	38, 51, 64, 79	0
19	S	150/154 (97%)	-0.19	0 100 100	32, 45, 66, 75	0
20	T	81/84 (96%)	-0.04	0 100 100	47, 65, 84, 89	0
21	U	119/119 (100%)	0.15	0 100 100	44, 62, 86, 97	0
22	V	53/66 (80%)	1.50	17 (32%) 1 0	85, 94, 102, 110	0
23	W	65/70 (92%)	0.42	3 (4%) 31 7	55, 81, 112, 118	0
24	X	154/154 (100%)	-0.20	0 100 100	32, 49, 66, 76	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
25	Y	82/91 (90%)	0.06	0 100 100	42, 58, 84, 99	0
26	Z	142/240 (59%)	-0.17	0 100 100	25, 46, 70, 85	0
27	1	73/73 (100%)	2.27	38 (52%) 0 0	79, 98, 103, 104	0
28	2	56/56 (100%)	-0.30	0 100 100	30, 39, 45, 49	0
29	3	46/48 (95%)	0.02	0 100 100	40, 66, 90, 102	0
30	4	92/92 (100%)	3.63	82 (89%) 0 0	91, 103, 108, 111	0
All	All	6577/7279 (90%)	-0.01	202 (3%) 47 9	23, 57, 102, 150	0

The worst 5 of 202 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
30	4	37	ASP	7.9
30	4	84	ARG	7.9
30	4	38	ARG	7.4
2	B	3001	U	6.7
30	4	1	MET	6.5

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

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

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
33	NA	A	8384	1/1	2.02	129.78	114,114,114,114	0
33	NA	A	8306	1/1	0.38	78.38	56,56,56,56	0
33	NA	A	8360	1/1	0.88	67.03	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	A	8356	1/1	0.69	43.27	58,58,58,58	0
33	NA	A	8329	1/1	0.48	36.99	70,70,70,70	0
32	MG	A	8024	1/1	0.67	35.90	116,116,116,116	0
34	CL	A	8515	1/1	0.58	25.80	100,100,100,100	0
33	NA	A	8363	1/1	0.40	25.01	66,66,66,66	0
33	NA	A	8385	1/1	0.32	24.21	41,41,41,41	0
33	NA	A	8382	1/1	0.44	22.56	62,62,62,62	0
33	NA	A	8372	1/1	0.68	22.25	55,55,55,55	0
32	MG	A	8070	1/1	0.60	21.48	66,66,66,66	0
33	NA	A	8316	1/1	0.32	21.23	51,51,51,51	0
33	NA	A	8370	1/1	0.30	19.29	49,49,49,49	0
33	NA	A	8364	1/1	0.27	19.27	40,40,40,40	0
33	NA	A	8379	1/1	0.34	17.56	41,41,41,41	0
34	CL	D	8519	1/1	0.54	16.39	65,65,65,65	0
33	NA	A	8371	1/1	0.31	16.27	54,54,54,54	0
33	NA	A	8321	1/1	0.44	14.31	39,39,39,39	0
32	MG	A	8082	1/1	0.21	14.11	52,52,52,52	0
33	NA	A	8374	1/1	0.44	13.34	63,63,63,63	0
33	NA	B	8383	1/1	0.54	13.26	63,63,63,63	0
33	NA	A	8362	1/1	0.36	12.64	69,69,69,69	0
34	CL	A	8522	1/1	0.54	12.63	75,75,75,75	0
33	NA	A	8322	1/1	0.43	12.22	46,46,46,46	0
33	NA	A	8340	1/1	0.25	12.22	31,31,31,31	0
34	CL	A	8505	1/1	0.66	11.97	88,88,88,88	0
33	NA	A	8352	1/1	0.39	11.87	52,52,52,52	0
32	MG	A	8119	1/1	0.36	11.43	71,71,71,71	0
34	CL	A	8503	1/1	0.34	11.32	50,50,50,50	0
32	MG	A	8103	1/1	0.34	11.04	55,55,55,55	0
32	MG	A	8049	1/1	0.55	10.97	89,89,89,89	0
33	NA	A	8331	1/1	0.37	10.55	61,61,61,61	0
33	NA	A	8376	1/1	0.30	9.22	78,78,78,78	0
33	NA	A	8378	1/1	0.44	9.02	37,37,37,37	0
33	NA	A	8313	1/1	0.26	8.59	63,63,63,63	0
33	NA	T	8312	1/1	0.77	8.35	124,124,124,124	0
36	CD	P	8405	1/1	0.39	8.17	169,169,169,169	0
33	NA	A	8323	1/1	0.25	8.14	50,50,50,50	0
31	SPR	A	9001	59/59	0.39	8.02	78,88,95,95	0
33	NA	A	8308	1/1	0.25	7.99	69,69,69,69	0
33	NA	S	8386	1/1	0.30	7.84	53,53,53,53	0
35	K	A	8603	1/1	0.34	7.66	88,88,88,88	0
33	NA	A	8373	1/1	0.45	7.52	59,59,59,59	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	A	8342	1/1	0.21	7.27	47,47,47,47	0
33	NA	A	8365	1/1	0.48	7.20	49,49,49,49	0
33	NA	A	8359	1/1	0.36	6.96	61,61,61,61	0
33	NA	A	8377	1/1	0.24	5.81	60,60,60,60	0
32	MG	A	8118	1/1	0.34	5.72	62,62,62,62	0
32	MG	A	8094	1/1	0.23	5.53	85,85,85,85	0
33	NA	A	8318	1/1	0.19	5.41	34,34,34,34	0
32	MG	A	8090	1/1	0.28	5.21	36,36,36,36	0
32	MG	A	8102	1/1	1.03	4.88	87,87,87,87	0
32	MG	A	8104	1/1	0.20	4.37	40,40,40,40	0
34	CL	C	8509	1/1	0.32	4.32	86,86,86,86	0
33	NA	S	8337	1/1	0.30	4.28	49,49,49,49	0
34	CL	R	8511	1/1	0.49	4.28	63,63,63,63	0
32	MG	A	8093	1/1	0.21	4.27	56,56,56,56	0
32	MG	A	8092	1/1	0.20	4.09	91,91,91,91	0
33	NA	A	8332	1/1	0.27	3.95	58,58,58,58	0
32	MG	Z	8109	1/1	0.23	3.74	53,53,53,53	0
33	NA	A	8366	1/1	0.26	3.31	49,49,49,49	0
32	MG	A	8097	1/1	0.19	3.08	44,44,44,44	0
33	NA	A	8350	1/1	0.16	2.99	34,34,34,34	0
34	CL	A	8517	1/1	0.25	2.83	55,55,55,55	0
33	NA	A	8355	1/1	0.37	2.81	55,55,55,55	0
33	NA	A	8328	1/1	0.18	2.80	45,45,45,45	0
33	NA	A	8367	1/1	0.19	2.29	52,52,52,52	0
32	MG	A	8101	1/1	0.18	2.28	55,55,55,55	0
33	NA	A	8336	1/1	0.20	2.09	49,49,49,49	0
33	NA	A	8368	1/1	0.16	1.91	47,47,47,47	0
32	MG	A	8110	1/1	0.16	1.86	47,47,47,47	0
32	MG	A	8089	1/1	0.18	1.86	84,84,84,84	0
34	CL	M	8510	1/1	0.45	1.67	87,87,87,87	0
32	MG	A	8076	1/1	0.19	1.62	71,71,71,71	0
35	K	A	8601	1/1	0.17	1.60	73,73,73,73	0
33	NA	A	8354	1/1	0.19	1.59	40,40,40,40	0
34	CL	A	8514	1/1	0.22	1.59	61,61,61,61	0
34	CL	Z	8520	1/1	0.17	1.41	35,35,35,35	0
32	MG	A	8067	1/1	0.21	1.32	50,50,50,50	0
34	CL	A	8512	1/1	0.19	1.27	32,32,32,32	0
34	CL	A	8516	1/1	0.21	1.13	44,44,44,44	0
32	MG	A	8081	1/1	0.19	1.04	58,58,58,58	0
32	MG	A	8114	1/1	0.51	0.97	92,92,92,92	0
33	NA	A	8303	1/1	0.17	0.95	51,51,51,51	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	A	8375	1/1	0.33	0.88	53,53,53,53	0
33	NA	A	8301	1/1	0.17	0.87	43,43,43,43	0
33	NA	A	8315	1/1	0.15	0.67	30,30,30,30	0
32	MG	A	8040	1/1	0.18	0.58	78,78,78,78	0
32	MG	A	8047	1/1	0.17	0.51	62,62,62,62	0
33	NA	A	8314	1/1	0.17	0.47	33,33,33,33	0
33	NA	A	8341	1/1	0.17	0.43	43,43,43,43	0
32	MG	A	8042	1/1	0.14	0.28	44,44,44,44	0
33	NA	A	8326	1/1	0.20	0.04	46,46,46,46	0
35	K	A	8602	1/1	0.23	0.01	68,68,68,68	0
32	MG	A	8116	1/1	0.17	-0.02	67,67,67,67	0
32	MG	A	8098	1/1	0.17	-0.04	50,50,50,50	0
33	NA	A	8369	1/1	0.33	-0.09	52,52,52,52	0
33	NA	A	8335	1/1	0.15	-0.09	52,52,52,52	0
32	MG	A	8044	1/1	0.14	-0.12	52,52,52,52	0
32	MG	A	8064	1/1	0.15	-0.12	24,24,24,24	0
34	CL	N	8518	1/1	0.21	-0.16	56,56,56,56	0
32	MG	1	8105	1/1	0.33	-0.18	44,44,44,44	0
33	NA	A	8361	1/1	0.13	-0.21	53,53,53,53	0
33	NA	A	8324	1/1	0.14	-0.27	51,51,51,51	0
34	CL	P	8508	1/1	0.17	-0.32	93,93,93,93	0
34	CL	K	8521	1/1	0.15	-0.35	46,46,46,46	0
34	CL	O	8507	1/1	0.25	-0.53	62,62,62,62	0
36	CD	4	8404	1/1	0.71	-0.56	156,156,156,156	0
33	NA	A	8333	1/1	0.15	-0.60	33,33,33,33	0
36	CD	V	8401	1/1	0.44	-0.67	142,142,142,142	0
33	NA	B	8351	1/1	0.25	-0.72	69,69,69,69	0
32	MG	A	8100	1/1	0.12	-0.76	69,69,69,69	0
33	NA	A	8330	1/1	0.16	-0.82	43,43,43,43	0
33	NA	J	8309	1/1	0.13	-0.85	21,21,21,21	0
32	MG	A	8088	1/1	0.15	-0.88	45,45,45,45	0
32	MG	A	8079	1/1	0.14	-0.90	39,39,39,39	0
33	NA	A	8310	1/1	0.14	-0.92	29,29,29,29	0
34	CL	S	8506	1/1	0.15	-0.98	46,46,46,46	0
32	MG	A	8012	1/1	0.13	-0.99	52,52,52,52	0
33	NA	A	8305	1/1	0.14	-1.10	34,34,34,34	0
32	MG	A	8106	1/1	0.09	-1.17	47,47,47,47	0
33	NA	A	8381	1/1	0.12	-1.22	51,51,51,51	0
32	MG	A	8066	1/1	0.12	-1.26	83,83,83,83	0
34	CL	K	8501	1/1	0.14	-1.31	56,56,56,56	0
34	CL	4	8504	1/1	0.61	-1.31	95,95,95,95	0
33	NA	M	8380	1/1	0.14	-1.34	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	E	8304	1/1	0.11	-1.41	35,35,35,35	0
32	MG	A	8051	1/1	0.12	-1.42	56,56,56,56	0
32	MG	A	8086	1/1	0.07	-1.44	50,50,50,50	0
32	MG	A	8071	1/1	0.14	-1.47	91,91,91,91	0
32	MG	A	8074	1/1	0.07	-1.64	31,31,31,31	0
34	CL	A	8513	1/1	0.12	-1.76	56,56,56,56	0
33	NA	A	8339	1/1	0.14	-1.77	16,16,16,16	0
33	NA	A	8334	1/1	0.07	-1.79	36,36,36,36	0
33	NA	A	8319	1/1	0.09	-1.90	52,52,52,52	0
33	NA	A	8320	1/1	0.13	-1.91	33,33,33,33	0
33	NA	A	8349	1/1	0.13	-1.94	53,53,53,53	0
33	NA	A	8353	1/1	0.11	-1.95	38,38,38,38	0
33	NA	A	8317	1/1	0.10	-2.07	27,27,27,27	0
36	CD	1	8403	1/1	0.28	-2.11	138,138,138,138	0
32	MG	A	8022	1/1	0.11	-2.23	41,41,41,41	0
32	MG	A	8018	1/1	0.10	-2.26	61,61,61,61	0
32	MG	A	8107	1/1	0.05	-2.34	47,47,47,47	0
32	MG	U	8073	1/1	0.14	-2.42	42,42,42,42	0
33	NA	A	8327	1/1	0.12	-2.50	32,32,32,32	0
32	MG	A	8062	1/1	0.11	-2.65	72,72,72,72	0
33	NA	K	8346	1/1	0.08	-2.66	27,27,27,27	0
32	MG	C	8065	1/1	0.10	-2.71	57,57,57,57	0
36	CD	2	8402	1/1	0.05	-2.74	59,59,59,59	0
33	NA	A	8338	1/1	0.07	-2.76	67,67,67,67	0
32	MG	A	8058	1/1	0.09	-2.84	43,43,43,43	0
32	MG	A	8112	1/1	0.13	-2.95	44,44,44,44	0
33	NA	R	8348	1/1	0.09	-3.01	37,37,37,37	0
32	MG	A	8015	1/1	0.09	-3.01	57,57,57,57	0
32	MG	A	8027	1/1	0.05	-3.05	63,63,63,63	0
32	MG	A	8001	1/1	0.12	-3.09	39,39,39,39	0
34	CL	K	8502	1/1	0.07	-3.11	52,52,52,52	0
32	MG	A	8055	1/1	0.08	-3.18	71,71,71,71	0
32	MG	A	8041	1/1	0.08	-3.20	46,46,46,46	0
32	MG	A	8060	1/1	0.10	-3.27	45,45,45,45	0
32	MG	A	8002	1/1	0.10	-3.52	31,31,31,31	0
32	MG	A	8085	1/1	0.13	-3.53	72,72,72,72	0
32	MG	A	8056	1/1	0.09	-3.55	53,53,53,53	0
32	MG	A	8111	1/1	0.08	-3.55	69,69,69,69	0
32	MG	A	8096	1/1	0.11	-3.57	53,53,53,53	0
32	MG	A	8108	1/1	0.07	-3.64	88,88,88,88	0
32	MG	A	8117	1/1	0.12	-3.70	31,31,31,31	0
32	MG	A	8046	1/1	0.07	-3.70	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
32	MG	A	8057	1/1	0.10	-3.75	49,49,49,49	0
33	NA	A	8343	1/1	0.09	-3.82	16,16,16,16	0
32	MG	A	8005	1/1	0.12	-3.83	44,44,44,44	0
32	MG	A	8068	1/1	0.07	-3.88	58,58,58,58	0
33	NA	A	8357	1/1	0.08	-3.89	67,67,67,67	0
32	MG	A	8030	1/1	0.10	-4.00	26,26,26,26	0
32	MG	A	8087	1/1	0.05	-4.24	48,48,48,48	0
32	MG	A	8038	1/1	0.08	-4.28	35,35,35,35	0
33	NA	A	8344	1/1	0.08	-4.34	30,30,30,30	0
33	NA	C	8345	1/1	0.10	-4.38	42,42,42,42	0
32	MG	A	8011	1/1	0.11	-4.44	52,52,52,52	0
33	NA	A	8307	1/1	0.10	-4.51	39,39,39,39	0
32	MG	A	8048	1/1	0.09	-4.64	45,45,45,45	0
33	NA	A	8311	1/1	0.09	-4.71	42,42,42,42	0
32	MG	A	8099	1/1	0.09	-4.80	38,38,38,38	0
32	MG	A	8091	1/1	0.08	-4.84	48,48,48,48	0
32	MG	A	8077	1/1	0.08	-4.85	31,31,31,31	0
32	MG	A	8004	1/1	0.07	-4.95	48,48,48,48	0
32	MG	A	8043	1/1	0.08	-5.04	39,39,39,39	0
32	MG	A	8006	1/1	0.08	-5.06	48,48,48,48	0
32	MG	L	8069	1/1	0.06	-5.18	50,50,50,50	0
32	MG	A	8045	1/1	0.10	-5.25	54,54,54,54	0
32	MG	A	8059	1/1	0.08	-5.39	31,31,31,31	0
32	MG	A	8039	1/1	0.06	-5.41	50,50,50,50	0
32	MG	A	8053	1/1	0.10	-5.58	52,52,52,52	0
32	MG	A	8033	1/1	0.07	-5.80	30,30,30,30	0
33	NA	A	8302	1/1	0.11	-5.81	40,40,40,40	0
32	MG	4	8078	1/1	0.31	-5.87	74,74,74,74	0
32	MG	A	8054	1/1	0.09	-6.01	48,48,48,48	0
33	NA	N	8347	1/1	0.07	-6.02	21,21,21,21	0
32	MG	A	8032	1/1	0.04	-6.10	34,34,34,34	0
32	MG	A	8009	1/1	0.05	-6.15	20,20,20,20	0
32	MG	A	8037	1/1	0.10	-6.23	48,48,48,48	0
32	MG	A	8003	1/1	0.08	-6.48	24,24,24,24	0
32	MG	A	8008	1/1	0.07	-6.54	49,49,49,49	0
32	MG	A	8017	1/1	0.04	-6.55	27,27,27,27	0
32	MG	A	8021	1/1	0.07	-6.76	27,27,27,27	0
32	MG	A	8014	1/1	0.07	-6.96	30,30,30,30	0
32	MG	A	8010	1/1	0.05	-7.08	40,40,40,40	0
32	MG	A	8035	1/1	0.05	-7.27	54,54,54,54	0
32	MG	A	8080	1/1	0.06	-7.28	50,50,50,50	0
32	MG	A	8026	1/1	0.04	-7.43	11,11,11,11	0

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*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
32	MG	A	8036	1/1	0.06	-7.45	45,45,45,45	0
32	MG	A	8013	1/1	0.11	-7.74	46,46,46,46	0
32	MG	A	8034	1/1	0.06	-7.87	39,39,39,39	0
32	MG	A	8020	1/1	0.05	-7.90	51,51,51,51	0
32	MG	A	8019	1/1	0.05	-7.96	35,35,35,35	0
32	MG	A	8063	1/1	0.06	-8.06	78,78,78,78	0
32	MG	A	8084	1/1	0.07	-8.28	48,48,48,48	0
32	MG	A	8061	1/1	0.08	-8.43	44,44,44,44	0
32	MG	A	8050	1/1	0.12	-8.70	85,85,85,85	0
32	MG	B	8095	1/1	0.08	-9.04	67,67,67,67	0
32	MG	A	8016	1/1	0.07	-9.05	41,41,41,41	0
32	MG	A	8072	1/1	0.13	-9.72	80,80,80,80	0
33	NA	A	8325	1/1	0.08	-10.10	52,52,52,52	0
32	MG	A	8023	1/1	0.07	-10.47	42,42,42,42	0
32	MG	A	8075	1/1	0.07	-10.90	57,57,57,57	0
32	MG	A	8115	1/1	0.10	-11.00	59,59,59,59	0
32	MG	A	8025	1/1	0.06	-13.85	60,60,60,60	0
32	MG	A	8007	1/1	0.04	-14.00	23,23,23,23	0
32	MG	A	8083	1/1	0.07	-14.90	47,47,47,47	0
32	MG	A	8029	1/1	0.06	-15.29	51,51,51,51	0
32	MG	A	8031	1/1	0.05	-18.82	31,31,31,31	0
32	MG	A	8028	1/1	0.07	-19.44	44,44,44,44	0
32	MG	A	8052	1/1	0.08	-20.58	45,45,45,45	0
32	MG	A	8113	1/1	0.12	-39.00	45,45,45,45	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.