



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 03:06 AM UTC

PDB ID : 1K73 / pdb_00001k73
Title : Co-crystal Structure of Anisomycin Bound to the 50S Ribosomal Subunit
Authors : Hansen, J.; Ban, N.; Nissen, P.; Moore, P.B.; Steitz, T.A.
Deposited on : 2001-10-18
Resolution : 3.01 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

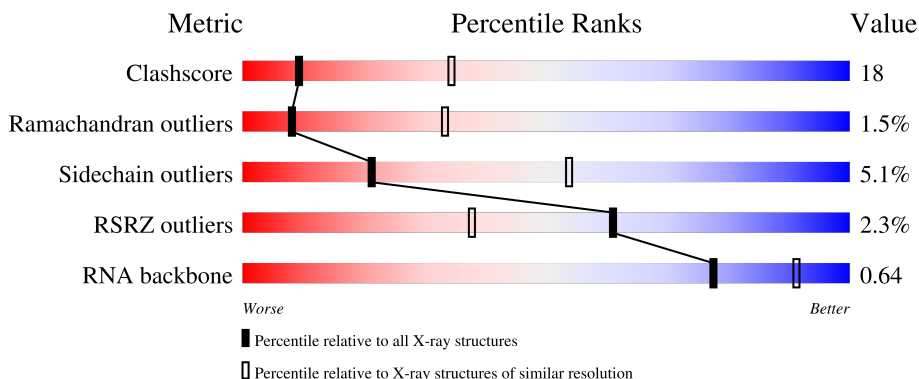
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

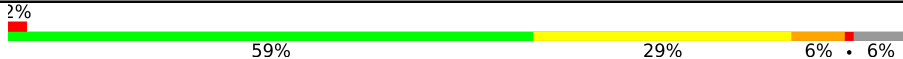
The reported resolution of this entry is 3.01 Å.

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.



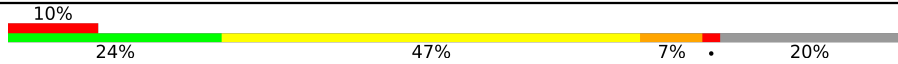
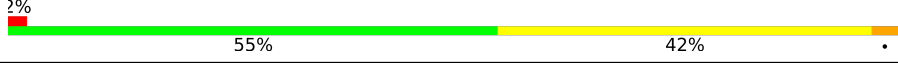
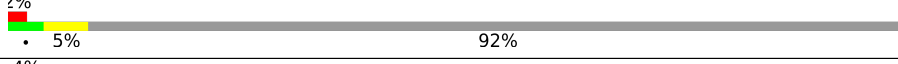
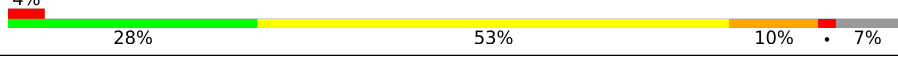
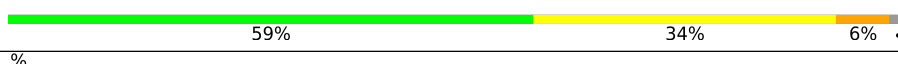
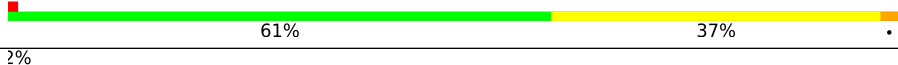
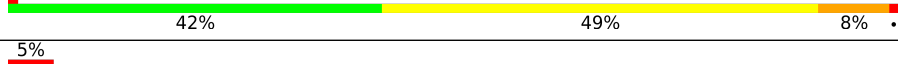
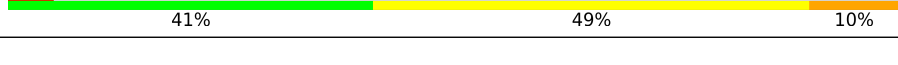
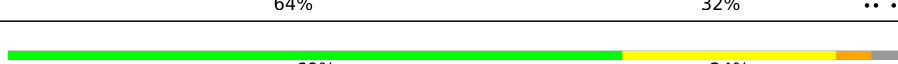
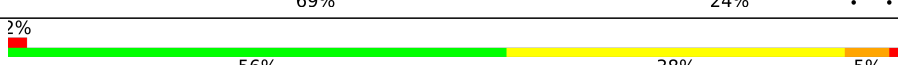
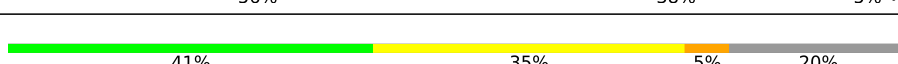
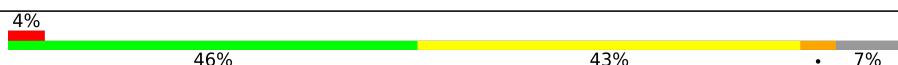
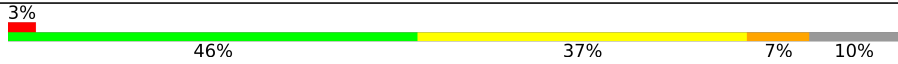
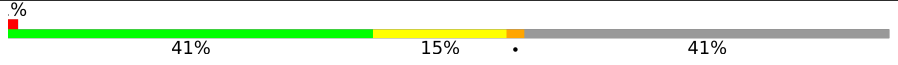
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)
RNA backbone	3983	1139 (3.24-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	2922	
2	B	122	
3	C	239	
4	D	337	
5	E	246	

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Mol	Chain	Length	Quality of chain
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	
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	

2 Entry composition

There are 37 unique types of molecules in this entry. The entry contains 98548 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	S	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	S	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	S	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			

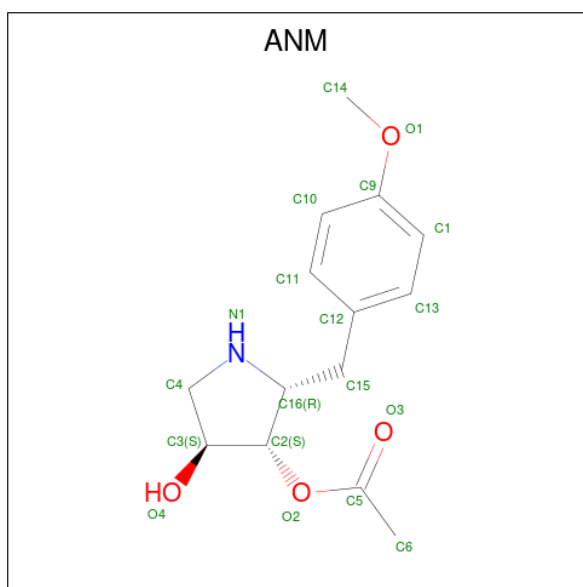
There is a discrepancy between the modelled and reference sequences:

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 ANISOMYCIN (CCD ID: ANM) (formula: C₁₄H₁₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
31	A	1	Total	C	N	O	0	0
			19	14	1	4		

- Molecule 32 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
32	A	109	Total	Mg	0	0
			109	109		
32	B	1	Total	Mg	0	0
			1	1		
32	C	2	Total	Mg	0	0
			2	2		
32	D	1	Total	Mg	0	0
			1	1		
32	L	1	Total	Mg	0	0
			1	1		
32	U	1	Total	Mg	0	0
			1	1		
32	Z	1	Total	Mg	0	0
			1	1		
32	4	1	Total	Mg	0	0
			1	1		

- Molecule 33 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
33	A	2	Total K 2 2	0	0

- Molecule 34 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
34	A	71	Total Na 71 71	0	0
34	B	2	Total Na 2 2	0	0
34	C	1	Total Na 1 1	0	0
34	E	1	Total Na 1 1	0	0
34	J	2	Total Na 2 2	0	0
34	K	1	Total Na 1 1	0	0
34	M	1	Total Na 1 1	0	0
34	N	1	Total Na 1 1	0	0
34	R	1	Total Na 1 1	0	0
34	S	3	Total Na 3 3	0	0
34	T	1	Total Na 1 1	0	0
34	U	1	Total Na 1 1	0	0

- Molecule 35 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
35	A	9	Total Cl 9 9	0	0
35	C	1	Total Cl 1 1	0	0
35	D	1	Total Cl 1 1	0	0
35	K	3	Total Cl 3 3	0	0
35	M	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
35	N	1	Total 1	Cl 1	0	0
35	O	1	Total 1	Cl 1	0	0
35	P	1	Total 1	Cl 1	0	0
35	R	1	Total 1	Cl 1	0	0
35	S	1	Total 1	Cl 1	0	0
35	Z	1	Total 1	Cl 1	0	0
35	4	1	Total 1	Cl 1	0	0

- Molecule 36 is CADMIUM ION (CCD ID: CD) (formula: Cd).

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

- Molecule 37 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	A	5924	Total 5924	O 5924	0	0
37	B	143	Total 143	O 143	0	0
37	C	127	Total 127	O 127	0	0
37	D	146	Total 146	O 146	0	0
37	E	164	Total 164	O 164	0	0

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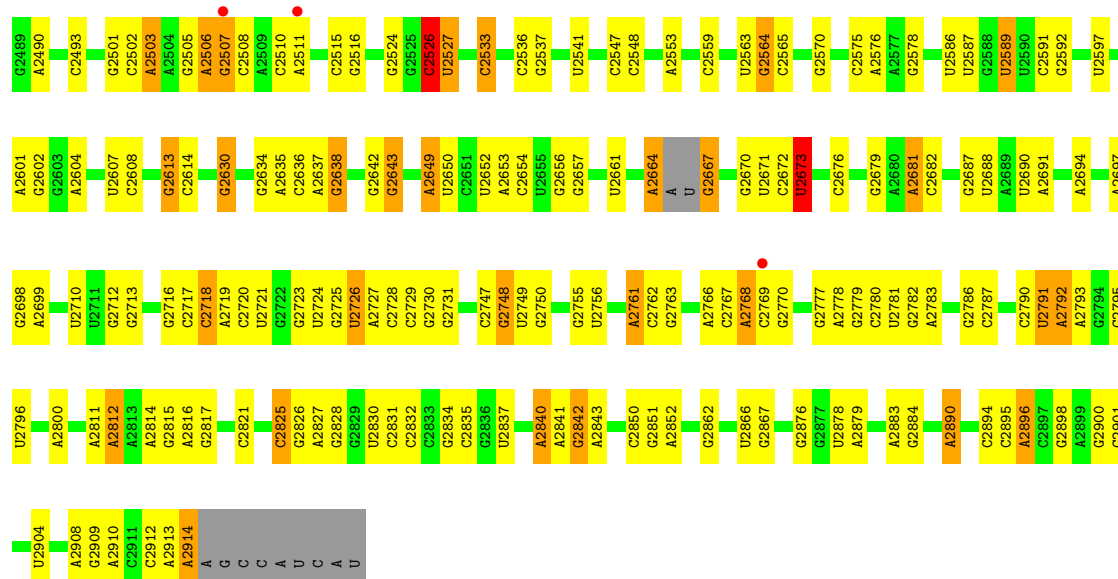
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	F	54	Total 54	O 54	0	0
37	G	42	Total 42	O 42	0	0
37	H	28	Total 28	O 28	0	0
37	I	22	Total 22	O 22	0	0
37	J	79	Total 79	O 79	0	0
37	K	56	Total 56	O 56	0	0
37	L	62	Total 62	O 62	0	0
37	M	81	Total 81	O 81	0	0
37	N	126	Total 126	O 126	0	0
37	O	66	Total 66	O 66	0	0
37	P	44	Total 44	O 44	0	0
37	Q	65	Total 65	O 65	0	0
37	R	54	Total 54	O 54	0	0
37	S	86	Total 86	O 86	0	0
37	T	35	Total 35	O 35	0	0
37	U	41	Total 41	O 41	0	0
37	V	25	Total 25	O 25	0	0
37	W	15	Total 15	O 15	0	0
37	X	67	Total 67	O 67	0	0
37	Y	29	Total 29	O 29	0	0
37	Z	92	Total 92	O 92	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	1	38	Total 38	O 38	0	0
37	2	56	Total 56	O 56	0	0
37	3	40	Total 40	O 40	0	0
37	4	72	Total 72	O 72	0	0

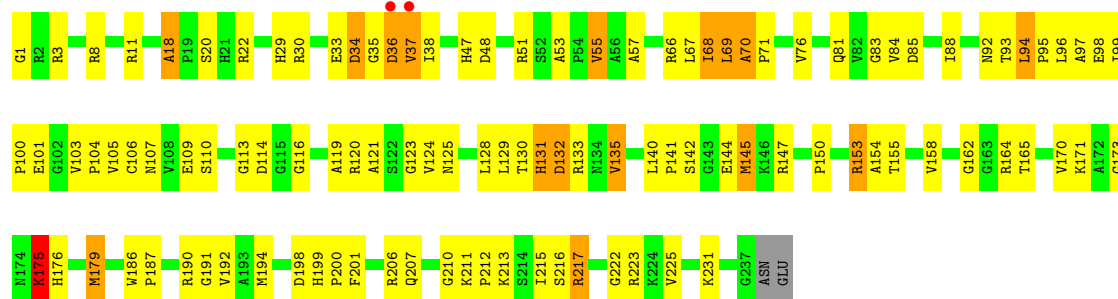




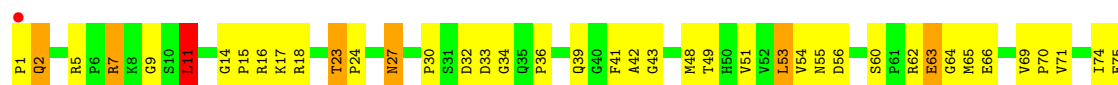
• Molecule 2: 5S RRNA

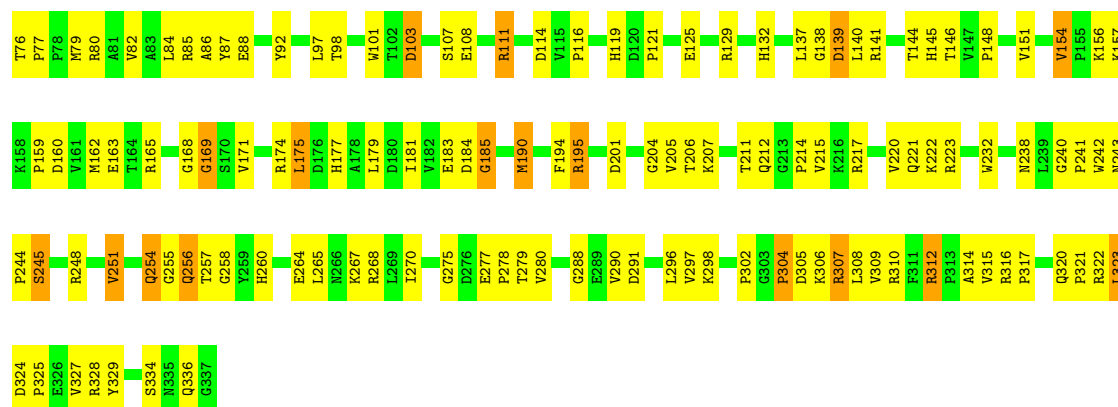


• Molecule 3: RIBOSOMAL PROTEIN L2



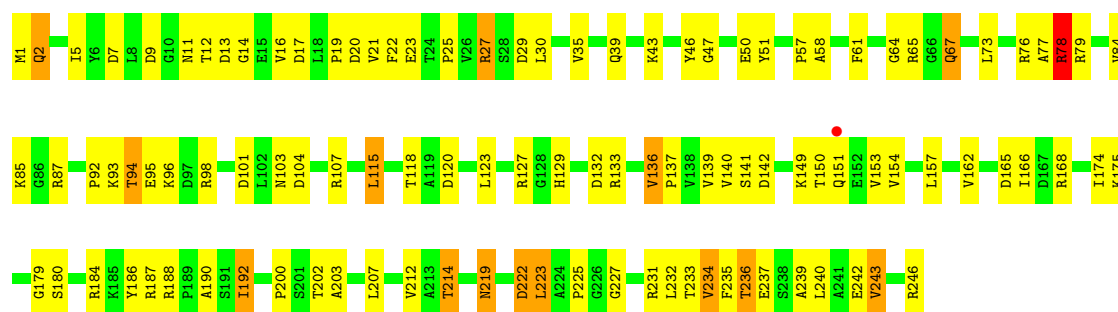
• Molecule 4: RIBOSOMAL PROTEIN L3





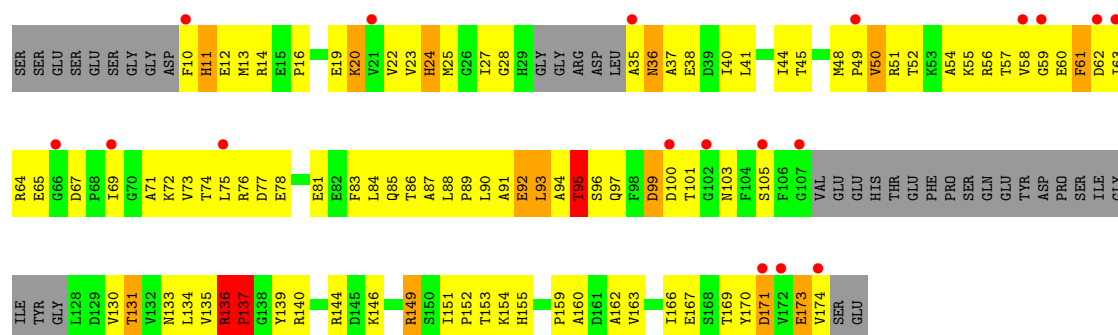
• Molecule 5: RIBOSOMAL PROTEIN L4

Chain E: 56% 38% 6%



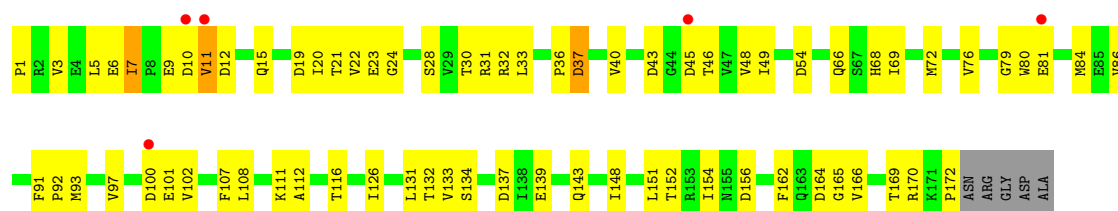
• Molecule 6: RIBOSOMAL PROTEIN L5

Chain F: 10% 24% 47% 7% 20%

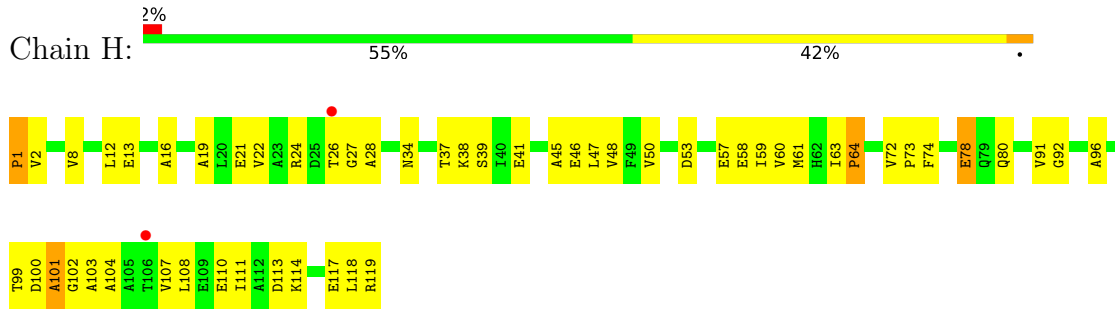


• Molecule 7: RIBOSOMAL PROTEIN L6

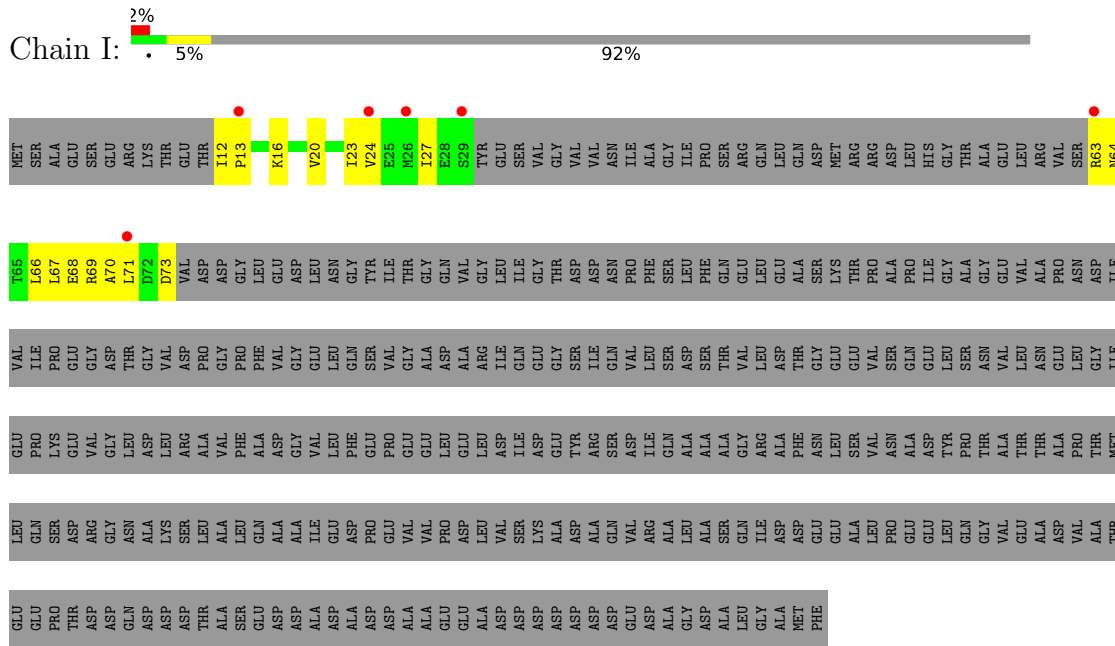
Chain G: 3% 56% 39% 2%



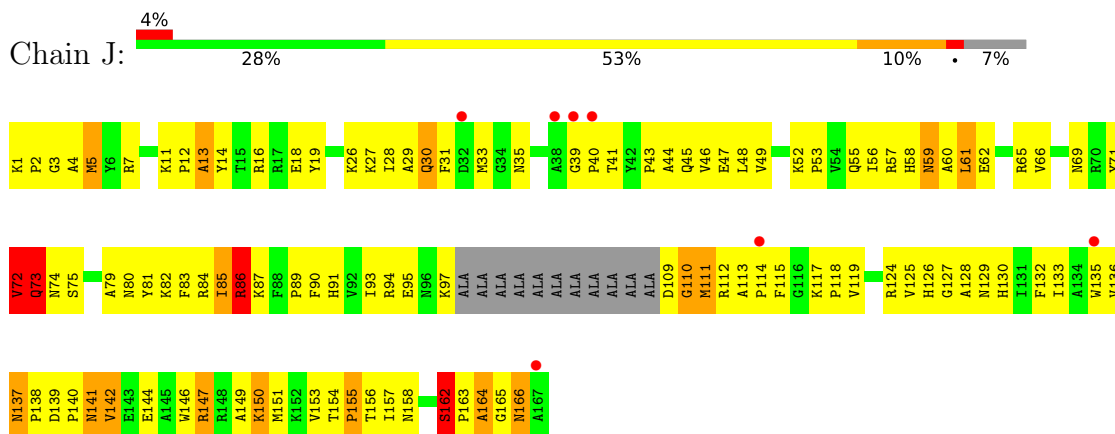
- Molecule 8: RIBOSOMAL PROTEIN L7AE



- Molecule 9: RIBOSOMAL PROTEIN L10

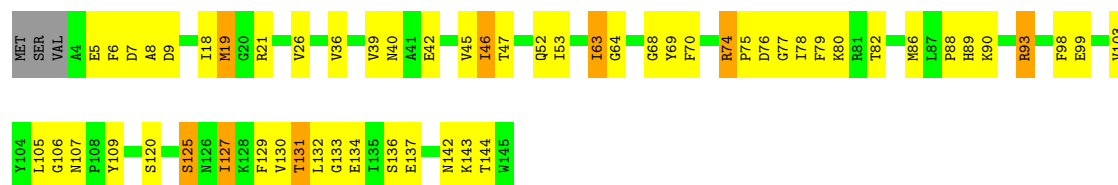


- Molecule 10: RIBOSOMAL PROTEIN L10E

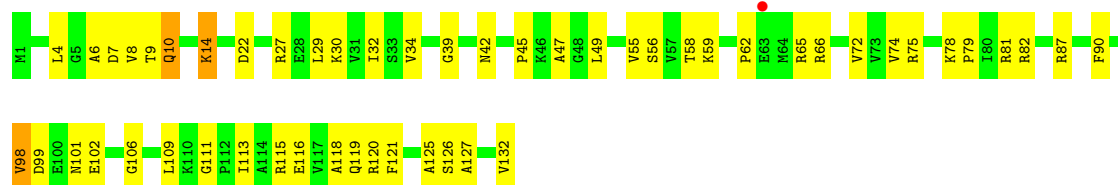


- Molecule 11: RIBOSOMAL PROTEIN L13

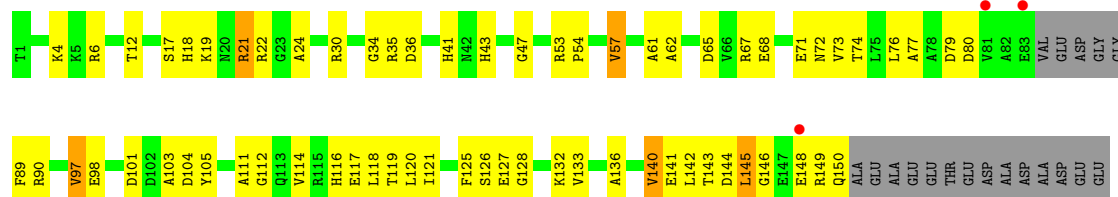




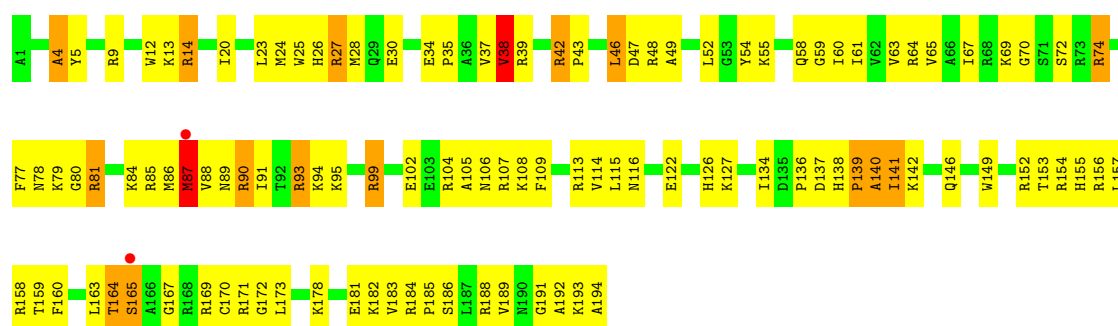
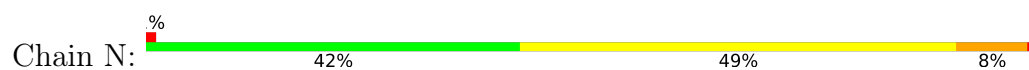
• Molecule 12: RIBOSOMAL PROTEIN L14



• Molecule 13: RIBOSOMAL PROTEIN L15

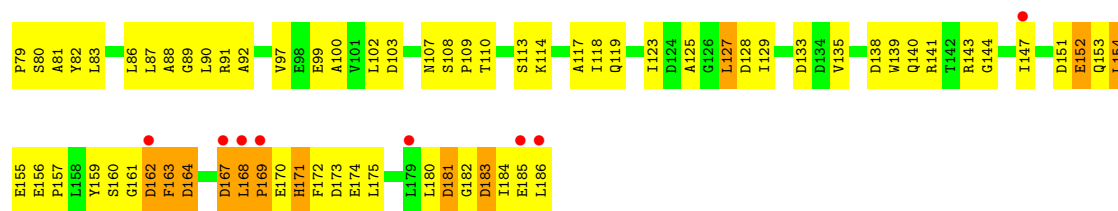


• Molecule 14: RIBOSOMAL PROTEIN L15E



• Molecule 15: RIBOSOMAL PROTEIN L18

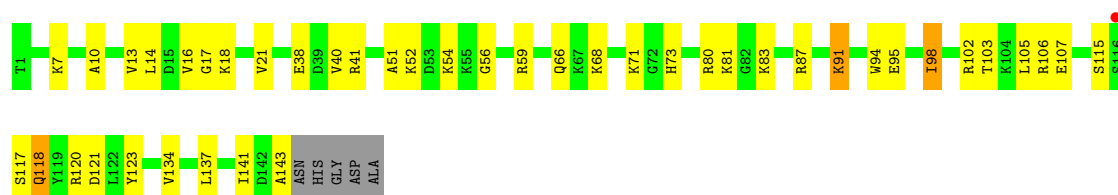




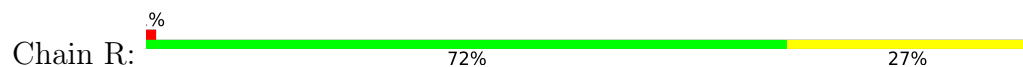
• Molecule 16: RIBOSOMAL PROTEIN L18E



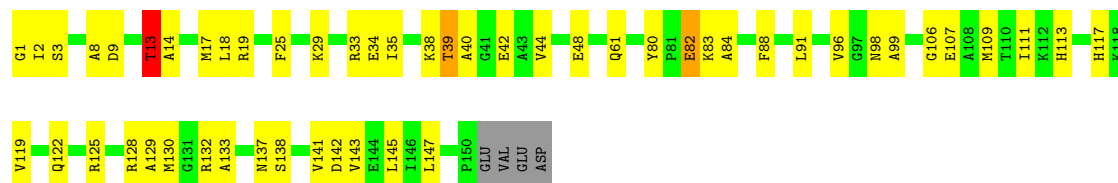
• Molecule 17: RIBOSOMAL PROTEIN L19E



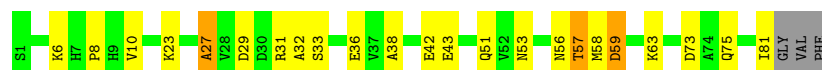
• Molecule 18: RIBOSOMAL PROTEIN L21E



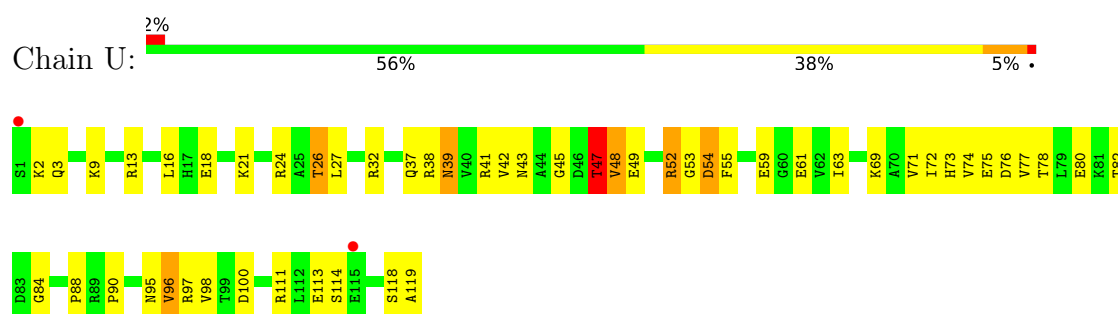
• Molecule 19: RIBOSOMAL PROTEIN L22



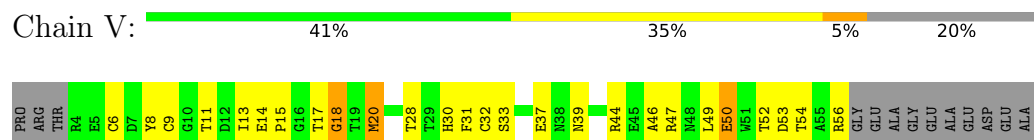
• Molecule 20: RIBOSOMAL PROTEIN L23



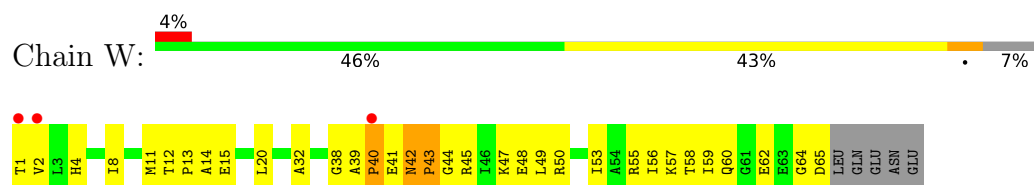
• Molecule 21: RIBOSOMAL PROTEIN L24



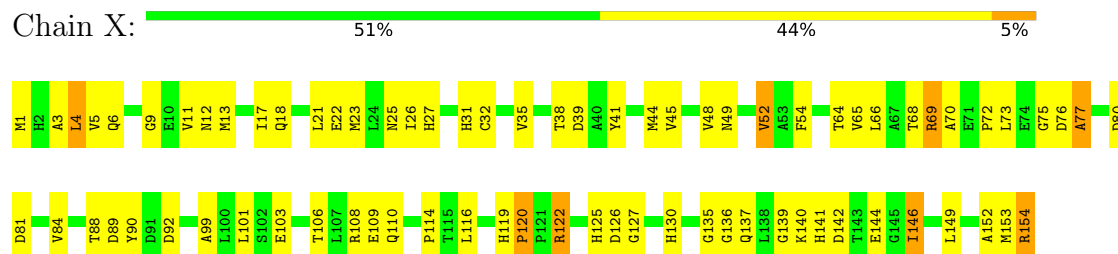
• Molecule 22: RIBOSOMAL PROTEIN L24E



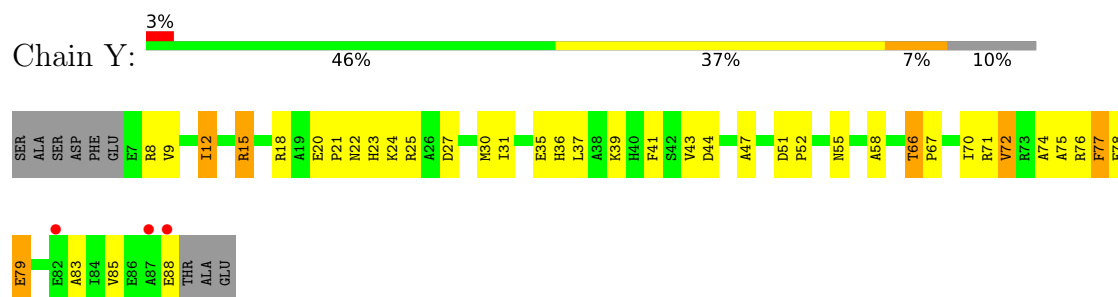
• Molecule 23: RIBOSOMAL PROTEIN L29



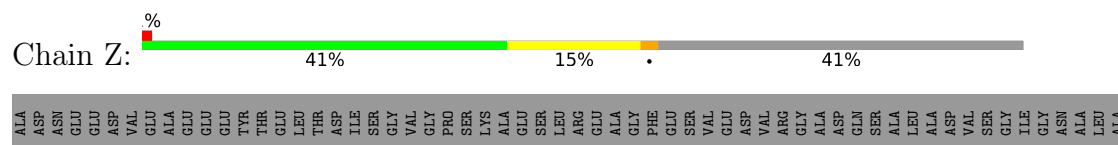
• Molecule 24: RIBOSOMAL PROTEIN L30

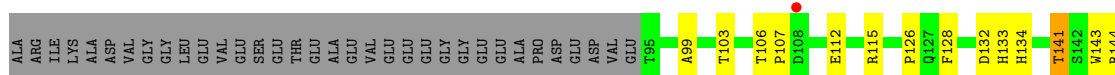


• Molecule 25: RIBOSOMAL PROTEIN L31E

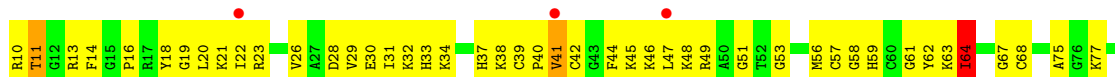


• Molecule 26: RIBOSOMAL PROTEIN L32E

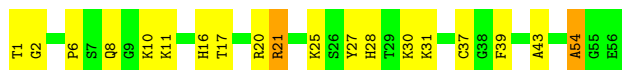




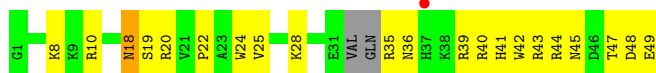
• Molecule 27: RIBOSOMAL PROTEIN L37Ae



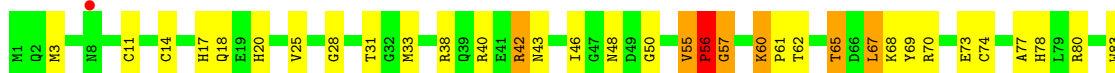
• Molecule 28: RIBOSOMAL PROTEIN L37E



• 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, α , β , γ	213.25Å 300.75Å 574.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.01 20.00 – 3.01	Depositor EDS
% Data completeness (in resolution range)	90.9 (20.00-3.01) 90.8 (20.00-3.01)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 3.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.212 , 0.246 0.209 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	98548	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CD, NA, ANM, MG, K, CL

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	0.57	12/66076 (0.0%)	0.77	61/103052 (0.1%)
2	B	0.55	1/2905 (0.0%)	0.83	6/4528 (0.1%)
3	C	0.68	1/1787 (0.1%)	1.11	7/2409 (0.3%)
4	D	0.68	0/2689	1.10	17/3652 (0.5%)
5	E	0.72	0/1883	1.12	10/2551 (0.4%)
6	F	0.54	0/1111	0.98	6/1498 (0.4%)
7	G	0.64	0/1382	0.97	5/1880 (0.3%)
8	H	0.64	0/896	0.97	2/1219 (0.2%)
9	I	0.55	0/241	0.83	0/324
10	J	0.77	2/1246 (0.2%)	1.33	16/1686 (0.9%)
11	K	0.72	2/1135 (0.2%)	1.02	3/1530 (0.2%)
12	L	0.64	0/1003	1.14	4/1351 (0.3%)
13	M	0.62	0/1126	1.13	6/1504 (0.4%)
14	N	0.74	1/1633 (0.1%)	1.18	13/2180 (0.6%)
15	O	0.65	0/1473	1.17	16/1999 (0.8%)
16	P	0.72	0/873	1.07	4/1181 (0.3%)
17	Q	0.61	0/1143	0.98	2/1521 (0.1%)
18	R	0.72	0/748	1.21	6/1005 (0.6%)
19	S	0.75	0/1172	1.13	6/1578 (0.4%)
20	T	0.62	0/648	0.98	3/875 (0.3%)
21	U	0.60	0/957	1.08	6/1289 (0.5%)
22	V	0.69	1/417 (0.2%)	1.03	2/562 (0.4%)
23	W	0.61	0/502	1.01	2/675 (0.3%)
24	X	0.76	0/1218	1.12	6/1655 (0.4%)
25	Y	0.71	0/664	1.06	3/895 (0.3%)
26	Z	0.67	0/1146	1.07	1/1536 (0.1%)
27	1	0.73	0/575	1.11	2/763 (0.3%)
28	2	0.72	0/437	1.09	3/578 (0.5%)
29	3	0.70	1/398 (0.3%)	0.84	0/527
30	4	0.64	0/771	1.06	5/1024 (0.5%)
All	All	0.60	21/98255 (0.0%)	0.87	223/147027 (0.2%)

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
1	A	1	58
2	B	1	2
All	All	2	60

The worst 5 of 21 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2486	A	O3'-P	32.86	2.10	1.61
1	A	2487	C	P-OP2	8.44	1.65	1.49
10	J	5	MET	SD-CE	-7.38	1.61	1.79
3	C	145	MET	SD-CE	6.95	1.97	1.79
11	K	19	MET	SD-CE	-6.57	1.63	1.79

The worst 5 of 223 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1164	U	OP1-P-O3'	-14.98	63.06	108.00
10	J	74	ASN	N-CA-C	-14.97	92.69	114.39
1	A	1563	G	C2'-C3'-O3'	13.94	130.40	109.50
2	B	3024	U	C2'-C3'-O3'	13.26	129.39	109.50
10	J	141	ASN	N-CA-C	-12.86	97.57	113.38

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	1563	G	C3'
2	B	3024	U	C3'

5 of 60 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	174	A	Sidechain
1	A	246	G	Sidechain
1	A	321	A	Sidechain
1	A	333	G	Sidechain
1	A	396	U	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	59017	0	29807	900	0
2	B	2600	0	1326	84	0
3	C	1754	0	1763	127	0
4	D	2624	0	2533	181	0
5	E	1858	0	1816	113	0
6	F	1094	0	1085	142	0
7	G	1357	0	1266	80	0
8	H	885	0	854	65	0
9	I	240	0	231	24	0
10	J	1215	0	1215	160	0
11	K	1119	0	1098	63	0
12	L	993	0	1027	56	0
13	M	1114	0	1072	63	0
14	N	1605	0	1676	154	0
15	O	1444	0	1401	124	0
16	P	864	0	873	30	0
17	Q	1133	0	1127	47	0
18	R	734	0	728	16	0
19	S	1149	0	1122	57	0
20	T	641	0	605	25	0
21	U	949	0	923	54	0
22	V	410	0	364	33	0
23	W	499	0	511	30	0
24	X	1195	0	1137	95	0
25	Y	654	0	653	46	0
26	Z	1130	0	1133	57	0
27	1	563	0	597	57	0
28	2	430	0	426	23	0
29	3	393	0	406	28	0
30	4	755	0	730	37	0
31	A	19	0	19	1	0
32	4	1	0	0	0	0
32	A	109	0	0	0	0
32	B	1	0	0	0	0
32	C	2	0	0	0	0
32	D	1	0	0	0	0
32	L	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	U	1	0	0	0	0
32	Z	1	0	0	0	0
33	A	2	0	0	0	0
34	A	71	0	0	0	0
34	B	2	0	0	0	0
34	C	1	0	0	0	0
34	E	1	0	0	0	0
34	J	2	0	0	0	0
34	K	1	0	0	0	0
34	M	1	0	0	0	0
34	N	1	0	0	0	0
34	R	1	0	0	0	0
34	S	3	0	0	0	0
34	T	1	0	0	0	0
34	U	1	0	0	0	0
35	4	1	0	0	0	0
35	A	9	0	0	0	0
35	C	1	0	0	0	0
35	D	1	0	0	0	0
35	K	3	0	0	2	0
35	M	1	0	0	0	0
35	N	1	0	0	1	0
35	O	1	0	0	0	0
35	P	1	0	0	0	0
35	R	1	0	0	0	0
35	S	1	0	0	0	0
35	Z	1	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	38	0	0	13	0
37	2	56	0	0	1	0
37	3	40	0	0	6	0
37	4	72	0	0	9	0
37	A	5924	0	0	183	0
37	B	143	0	0	15	0
37	C	127	0	0	23	0
37	D	146	0	0	25	0
37	E	164	0	0	28	0
37	F	54	0	0	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
37	G	42	0	0	11	0
37	H	28	0	0	9	0
37	I	22	0	0	6	0
37	J	79	0	0	23	0
37	K	56	0	0	5	0
37	L	62	0	0	9	0
37	M	81	0	0	18	0
37	N	126	0	0	21	0
37	O	66	0	0	17	0
37	P	44	0	0	5	0
37	Q	65	0	0	3	0
37	R	54	0	0	2	0
37	S	86	0	0	10	0
37	T	35	0	0	5	0
37	U	41	0	0	4	0
37	V	25	0	0	6	0
37	W	15	0	0	2	0
37	X	67	0	0	8	0
37	Y	29	0	0	4	0
37	Z	92	0	0	16	0
All	All	98548	0	59524	2713	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

The worst 5 of 2713 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1160:G:H5'	1:A:1161:A:H5'	1.24	1.18
5:E:236:THR:HG22	5:E:239:ALA:H	1.04	1.14
10:J:86:ARG:NH1	10:J:133:ILE:HG13	1.64	1.12
1:A:1134:G:H4'	10:J:151:MET:HE1	1.26	1.11
6:F:25:MET:HE2	6:F:41:LEU:HG	1.31	1.10

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%)	212 (90%)	19 (8%)	4 (2%)	7	30
4	D	335/337 (99%)	310 (92%)	18 (5%)	7 (2%)	5	25
5	E	244/246 (99%)	222 (91%)	22 (9%)	0	100	100
6	F	134/176 (76%)	99 (74%)	24 (18%)	11 (8%)	0	3
7	G	170/177 (96%)	162 (95%)	8 (5%)	0	100	100
8	H	117/119 (98%)	105 (90%)	10 (8%)	2 (2%)	7	30
9	I	25/348 (7%)	24 (96%)	1 (4%)	0	100	100
10	J	152/167 (91%)	132 (87%)	15 (10%)	5 (3%)	3	16
11	K	140/145 (97%)	130 (93%)	7 (5%)	3 (2%)	5	25
12	L	130/132 (98%)	121 (93%)	7 (5%)	2 (2%)	8	33
13	M	141/164 (86%)	122 (86%)	17 (12%)	2 (1%)	9	34
14	N	192/194 (99%)	177 (92%)	12 (6%)	3 (2%)	7	32
15	O	184/186 (99%)	167 (91%)	10 (5%)	7 (4%)	2	13
16	P	113/115 (98%)	109 (96%)	4 (4%)	0	100	100
17	Q	141/148 (95%)	137 (97%)	4 (3%)	0	100	100
18	R	93/95 (98%)	87 (94%)	6 (6%)	0	100	100
19	S	148/154 (96%)	142 (96%)	6 (4%)	0	100	100
20	T	79/84 (94%)	75 (95%)	4 (5%)	0	100	100
21	U	117/119 (98%)	111 (95%)	6 (5%)	0	100	100
22	V	51/66 (77%)	47 (92%)	4 (8%)	0	100	100
23	W	63/70 (90%)	58 (92%)	3 (5%)	2 (3%)	3	17
24	X	152/154 (99%)	146 (96%)	4 (3%)	2 (1%)	9	36
25	Y	80/91 (88%)	70 (88%)	9 (11%)	1 (1%)	9	36
26	Z	140/240 (58%)	140 (100%)	0	0	100	100
27	1	71/73 (97%)	64 (90%)	5 (7%)	2 (3%)	4	19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
28	2	54/56 (96%)	53 (98%)	1 (2%)	0	100	100
29	3	42/48 (88%)	42 (100%)	0	0	100	100
30	4	90/92 (98%)	87 (97%)	1 (1%)	2 (2%)	5	24
All	All	3633/4235 (86%)	3351 (92%)	227 (6%)	55 (2%)	8	33

5 of 55 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%)	168 (94%)	11 (6%)	17	47
4	D	282/282 (100%)	263 (93%)	19 (7%)	15	44
5	E	193/193 (100%)	177 (92%)	16 (8%)	10	35
6	F	117/147 (80%)	108 (92%)	9 (8%)	12	38
7	G	152/155 (98%)	147 (97%)	5 (3%)	33	65
8	H	92/92 (100%)	89 (97%)	3 (3%)	33	65
9	I	27/283 (10%)	27 (100%)	0	100	100
10	J	122/122 (100%)	109 (89%)	13 (11%)	6	25
11	K	118/121 (98%)	111 (94%)	7 (6%)	18	49
12	L	106/106 (100%)	103 (97%)	3 (3%)	38	69
13	M	112/126 (89%)	109 (97%)	3 (3%)	39	70
14	N	166/166 (100%)	159 (96%)	7 (4%)	26	59
15	O	149/149 (100%)	142 (95%)	7 (5%)	23	56

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
16	P	93/93 (100%)	89 (96%)	4 (4%)	26	59
17	Q	113/116 (97%)	110 (97%)	3 (3%)	39	70
18	R	79/79 (100%)	74 (94%)	5 (6%)	16	46
19	S	117/121 (97%)	114 (97%)	3 (3%)	40	71
20	T	71/73 (97%)	70 (99%)	1 (1%)	59	79
21	U	105/105 (100%)	100 (95%)	5 (5%)	23	56
22	V	44/52 (85%)	44 (100%)	0	100	100
23	W	51/56 (91%)	50 (98%)	1 (2%)	48	75
24	X	130/130 (100%)	121 (93%)	9 (7%)	14	43
25	Y	66/73 (90%)	63 (96%)	3 (4%)	24	57
26	Z	120/195 (62%)	111 (92%)	9 (8%)	12	39
27	1	56/56 (100%)	54 (96%)	2 (4%)	31	64
28	2	46/46 (100%)	46 (100%)	0	100	100
29	3	42/44 (96%)	41 (98%)	1 (2%)	43	72
30	4	79/79 (100%)	75 (95%)	4 (5%)	21	54
All	All	3027/3441 (88%)	2874 (95%)	153 (5%)	21	54

5 of 153 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
19	S	13	THR
26	Z	200	THR
21	U	26	THR
24	X	122	ARG
30	4	42	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 111 such sidechains are listed below:

Mol	Chain	Res	Type
15	O	107	ASN
30	4	48	ASN
19	S	113	HIS
30	4	30	GLN
27	1	70	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	2747/2922 (94%)	244 (8%)	35 (1%)
2	B	121/122 (99%)	19 (15%)	7 (5%)
All	All	2868/3044 (94%)	263 (9%)	42 (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 42 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	2467	A
2	B	3002	U
1	A	2526	C
1	A	2718	C
2	B	3023	U

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 233 ligands modelled in this entry, 232 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	ANM	A	9000	33	20,20,20	1.78	4 (20%)	24,27,27	1.71	7 (29%)

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	ANM	A	9000	33	-	2/10/23/23	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
31	A	9000	ANM	C1-C9	5.55	1.49	1.38
31	A	9000	ANM	C10-C9	2.92	1.44	1.38
31	A	9000	ANM	C11-C12	2.80	1.44	1.38
31	A	9000	ANM	O1-C9	2.60	1.42	1.37

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	A	9000	ANM	C2-O2-C5	3.49	123.15	117.72
31	A	9000	ANM	C10-C9-C1	-3.29	115.37	120.16
31	A	9000	ANM	C13-C1-C9	2.78	122.91	119.73
31	A	9000	ANM	C12-C15-C16	-2.55	109.08	113.40
31	A	9000	ANM	O2-C5-O3	2.12	127.09	122.99

There are no chirality outliers.

All (2) torsion outliers are listed below:

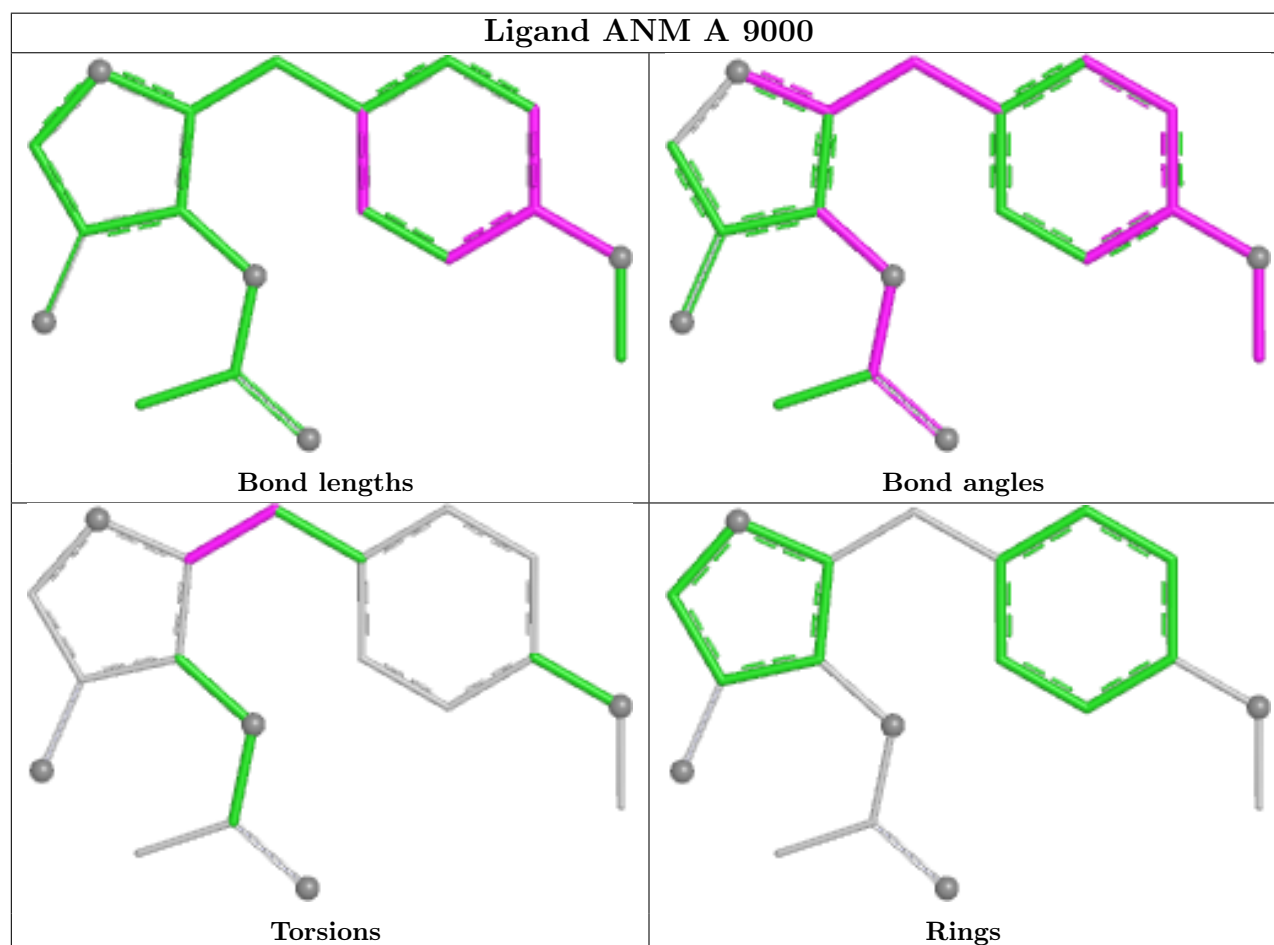
Mol	Chain	Res	Type	Atoms
31	A	9000	ANM	C12-C15-C16-C2
31	A	9000	ANM	C12-C15-C16-N1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
31	A	9000	ANM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	2486:A	O3'	2487:C	P	2.10

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	2754/2922 (94%)	-0.12	67 (2%) 59 36	20, 45, 89, 137	0
2	B	122/122 (100%)	0.48	7 (5%) 29 15	32, 61, 88, 146	0
3	C	237/239 (99%)	0.02	2 (0%) 82 64	27, 49, 83, 104	0
4	D	337/337 (100%)	0.09	1 (0%) 90 79	25, 54, 79, 88	0
5	E	246/246 (100%)	-0.00	1 (0%) 88 76	18, 44, 68, 79	0
6	F	140/176 (79%)	1.10	18 (12%) 7 4	51, 96, 114, 118	0
7	G	172/177 (97%)	0.46	5 (2%) 53 31	44, 66, 85, 91	0
8	H	119/119 (100%)	0.52	2 (1%) 69 46	50, 68, 93, 100	0
9	I	29/348 (8%)	1.34	6 (20%) 2 2	73, 88, 96, 100	0
10	J	156/167 (93%)	0.40	7 (4%) 38 20	34, 55, 84, 88	0
11	K	142/145 (97%)	0.03	0 100 100	34, 49, 70, 90	0
12	L	132/132 (100%)	0.06	1 (0%) 82 64	30, 50, 71, 78	0
13	M	145/164 (88%)	0.31	3 (2%) 63 40	23, 64, 101, 112	0
14	N	194/194 (100%)	0.04	2 (1%) 79 59	29, 43, 60, 72	0
15	O	186/186 (100%)	0.60	9 (4%) 35 18	40, 61, 101, 114	0
16	P	115/115 (100%)	0.23	0 100 100	38, 53, 69, 75	0
17	Q	143/148 (96%)	0.10	1 (0%) 84 66	35, 54, 68, 75	0
18	R	95/95 (100%)	-0.07	1 (1%) 78 57	33, 43, 58, 71	0
19	S	150/154 (97%)	-0.24	0 100 100	24, 43, 65, 71	0
20	T	81/84 (96%)	0.11	0 100 100	41, 58, 78, 81	0
21	U	119/119 (100%)	0.26	2 (1%) 69 46	37, 56, 79, 92	0
22	V	53/66 (80%)	0.29	0 100 100	40, 55, 72, 79	0
23	W	65/70 (92%)	0.48	3 (4%) 37 20	49, 71, 105, 112	0
24	X	154/154 (100%)	-0.08	0 100 100	34, 46, 65, 75	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
25	Y	82/91 (90%)	0.26	3 (3%) 45 25	42, 57, 82, 99	0
26	Z	142/240 (59%)	-0.06	2 (1%) 73 51	26, 44, 65, 83	0
27	1	73/73 (100%)	0.56	3 (4%) 41 22	45, 59, 75, 84	0
28	2	56/56 (100%)	-0.40	0 100 100	21, 34, 39, 41	0
29	3	46/48 (95%)	0.38	1 (2%) 62 39	35, 60, 85, 96	0
30	4	92/92 (100%)	0.19	1 (1%) 78 57	35, 55, 69, 80	0
All	All	6577/7279 (90%)	0.07	148 (2%) 61 38	18, 50, 90, 146	0

The worst 5 of 148 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1173	A	6.5
23	W	1	THR	6.4
2	B	3025	G	6.3
1	A	1175	G	5.4
1	A	1176	C	4.6

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 oligosaccharides 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. 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	RSCC	RSR	B-factors(Å ²)	Q<0.9
34	NA	A	8329	1/1	0.38	0.16	60,60,60,60	0
34	NA	A	8371	1/1	0.48	0.38	47,47,47,47	0
34	NA	S	8386	1/1	0.63	0.31	83,83,83,83	0
34	NA	A	8384	1/1	0.64	0.17	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
34	NA	A	8306	1/1	0.66	0.17	39,39,39,39	0
34	NA	A	8385	1/1	0.67	0.24	50,50,50,50	0
34	NA	A	8363	1/1	0.68	0.35	54,54,54,54	0
34	NA	B	8383	1/1	0.69	0.32	52,52,52,52	0
32	MG	A	8106	1/1	0.72	0.13	50,50,50,50	0
34	NA	A	8382	1/1	0.72	0.35	71,71,71,71	0
34	NA	A	8365	1/1	0.72	0.11	36,36,36,36	0
34	NA	B	8351	1/1	0.75	0.14	30,30,30,30	0
32	MG	A	8066	1/1	0.75	0.17	94,94,94,94	0
32	MG	A	8049	1/1	0.75	0.14	68,68,68,68	0
32	MG	A	8092	1/1	0.76	0.13	75,75,75,75	0
34	NA	A	8368	1/1	0.77	0.24	57,57,57,57	0
34	NA	A	8340	1/1	0.78	0.14	47,47,47,47	0
34	NA	A	8367	1/1	0.78	0.11	41,41,41,41	0
32	MG	A	8102	1/1	0.80	0.12	48,48,48,48	0
34	NA	A	8328	1/1	0.80	0.17	41,41,41,41	0
32	MG	A	8087	1/1	0.80	0.12	67,67,67,67	0
34	NA	T	8312	1/1	0.80	0.10	39,39,39,39	0
34	NA	A	8352	1/1	0.81	0.13	49,49,49,49	0
34	NA	A	8381	1/1	0.82	0.16	44,44,44,44	0
34	NA	A	8326	1/1	0.82	0.21	61,61,61,61	0
34	NA	A	8324	1/1	0.83	0.24	43,43,43,43	0
34	NA	A	8369	1/1	0.83	0.30	49,49,49,49	0
32	MG	A	8050	1/1	0.83	0.09	57,57,57,57	0
34	NA	A	8374	1/1	0.83	0.34	65,65,65,65	0
34	NA	A	8366	1/1	0.83	0.16	60,60,60,60	0
34	NA	A	8313	1/1	0.83	0.15	64,64,64,64	0
34	NA	A	8349	1/1	0.84	0.18	49,49,49,49	0
34	NA	A	8377	1/1	0.84	0.25	61,61,61,61	0
34	NA	A	8302	1/1	0.84	0.14	44,44,44,44	0
32	MG	B	8095	1/1	0.84	0.23	75,75,75,75	0
34	NA	A	8372	1/1	0.84	0.29	60,60,60,60	0
34	NA	A	8301	1/1	0.85	0.08	16,16,16,16	0
34	NA	A	8359	1/1	0.85	0.20	53,53,53,53	0
34	NA	A	8362	1/1	0.85	0.28	66,66,66,66	0
32	MG	U	8073	1/1	0.86	0.07	52,52,52,52	0
34	NA	A	8333	1/1	0.86	0.13	38,38,38,38	0
32	MG	A	8090	1/1	0.86	0.27	55,55,55,55	0
32	MG	A	8045	1/1	0.86	0.10	49,49,49,49	0
32	MG	A	8094	1/1	0.86	0.16	65,65,65,65	0
35	CL	A	8522	1/1	0.86	0.29	85,85,85,85	0
34	NA	A	8379	1/1	0.87	0.16	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
34	NA	A	8341	1/1	0.87	0.06	36,36,36,36	0
34	NA	A	8323	1/1	0.87	0.13	33,33,33,33	0
34	NA	A	8332	1/1	0.87	0.19	42,42,42,42	0
34	NA	A	8358	1/1	0.87	0.31	86,86,86,86	0
35	CL	Z	8520	1/1	0.87	0.14	39,39,39,39	0
35	CL	A	8517	1/1	0.88	0.08	52,52,52,52	0
32	MG	A	8088	1/1	0.88	0.12	24,24,24,24	0
34	NA	A	8376	1/1	0.88	0.17	38,38,38,38	0
34	NA	A	8357	1/1	0.89	0.16	50,50,50,50	0
34	NA	A	8319	1/1	0.89	0.07	23,23,23,23	0
32	MG	A	8081	1/1	0.89	0.07	60,60,60,60	0
34	NA	A	8361	1/1	0.89	0.15	60,60,60,60	0
34	NA	A	8373	1/1	0.89	0.21	49,49,49,49	0
34	NA	R	8348	1/1	0.89	0.04	34,34,34,34	0
34	NA	S	8337	1/1	0.89	0.09	41,41,41,41	0
32	MG	A	8091	1/1	0.89	0.07	52,52,52,52	0
32	MG	A	8051	1/1	0.89	0.10	70,70,70,70	0
34	NA	A	8311	1/1	0.89	0.10	55,55,55,55	0
32	MG	A	8076	1/1	0.89	0.11	68,68,68,68	0
35	CL	N	8518	1/1	0.89	0.09	53,53,53,53	0
34	NA	A	8355	1/1	0.89	0.20	58,58,58,58	0
34	NA	A	8330	1/1	0.90	0.07	41,41,41,41	0
34	NA	A	8331	1/1	0.90	0.14	51,51,51,51	0
34	NA	A	8308	1/1	0.90	0.14	53,53,53,53	0
34	NA	J	8322	1/1	0.90	0.23	59,59,59,59	0
35	CL	C	8509	1/1	0.90	0.10	64,64,64,64	0
35	CL	M	8510	1/1	0.90	0.10	74,74,74,74	0
32	MG	A	8114	1/1	0.90	0.07	54,54,54,54	0
35	CL	O	8507	1/1	0.90	0.08	65,65,65,65	0
32	MG	A	8024	1/1	0.90	0.07	29,29,29,29	0
32	MG	A	8052	1/1	0.91	0.09	63,63,63,63	0
32	MG	C	8105	1/1	0.91	0.12	6,6,6,6	0
32	MG	A	8046	1/1	0.91	0.07	53,53,53,53	0
32	MG	A	8103	1/1	0.91	0.13	49,49,49,49	0
34	NA	A	8354	1/1	0.91	0.17	39,39,39,39	0
35	CL	K	8502	1/1	0.91	0.07	62,62,62,62	0
34	NA	A	8375	1/1	0.91	0.16	47,47,47,47	0
34	NA	A	8321	1/1	0.91	0.18	39,39,39,39	0
32	MG	A	8082	1/1	0.91	0.14	60,60,60,60	0
32	MG	A	8067	1/1	0.91	0.12	48,48,48,48	0
35	CL	4	8504	1/1	0.91	0.10	54,54,54,54	0
31	ANM	A	9000	19/19	0.92	0.14	27,35,41,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
32	MG	A	8041	1/1	0.92	0.09	43,43,43,43	0
32	MG	A	8093	1/1	0.92	0.05	37,37,37,37	0
32	MG	A	8042	1/1	0.92	0.08	41,41,41,41	0
34	NA	A	8360	1/1	0.92	0.21	53,53,53,53	0
32	MG	A	8099	1/1	0.92	0.08	47,47,47,47	0
32	MG	A	8101	1/1	0.92	0.13	55,55,55,55	0
35	CL	R	8511	1/1	0.92	0.11	63,63,63,63	0
32	MG	A	8064	1/1	0.92	0.15	20,20,20,20	0
35	CL	A	8505	1/1	0.92	0.19	85,85,85,85	0
34	NA	A	8364	1/1	0.93	0.16	43,43,43,43	0
35	CL	A	8503	1/1	0.93	0.08	60,60,60,60	0
34	NA	A	8350	1/1	0.93	0.14	45,45,45,45	0
35	CL	A	8512	1/1	0.93	0.11	22,22,22,22	0
35	CL	A	8514	1/1	0.93	0.11	67,67,67,67	0
32	MG	A	8104	1/1	0.93	0.06	42,42,42,42	0
34	NA	A	8314	1/1	0.93	0.06	13,13,13,13	0
34	NA	A	8316	1/1	0.93	0.20	52,52,52,52	0
34	NA	A	8318	1/1	0.93	0.13	40,40,40,40	0
34	NA	A	8305	1/1	0.93	0.07	34,34,34,34	0
32	MG	A	8116	1/1	0.93	0.06	46,46,46,46	0
34	NA	A	8335	1/1	0.93	0.15	52,52,52,52	0
33	K	A	8201	1/1	0.93	0.15	69,69,69,69	0
34	NA	A	8310	1/1	0.93	0.10	32,32,32,32	0
32	MG	A	8111	1/1	0.93	0.06	49,49,49,49	0
32	MG	A	8089	1/1	0.94	0.11	63,63,63,63	0
34	NA	A	8378	1/1	0.94	0.15	41,41,41,41	0
32	MG	A	8075	1/1	0.94	0.06	51,51,51,51	0
32	MG	A	8096	1/1	0.94	0.04	52,52,52,52	0
32	MG	A	8053	1/1	0.94	0.07	34,34,34,34	0
34	NA	A	8356	1/1	0.94	0.17	53,53,53,53	0
32	MG	A	8010	1/1	0.94	0.03	26,26,26,26	0
34	NA	A	8325	1/1	0.94	0.09	51,51,51,51	0
34	NA	A	8336	1/1	0.94	0.05	45,45,45,45	0
34	NA	E	8304	1/1	0.94	0.09	35,35,35,35	0
34	NA	A	8339	1/1	0.94	0.04	15,15,15,15	0
34	NA	K	8346	1/1	0.94	0.07	35,35,35,35	0
34	NA	A	8315	1/1	0.94	0.12	36,36,36,36	0
34	NA	A	8327	1/1	0.94	0.10	18,18,18,18	0
34	NA	A	8307	1/1	0.94	0.07	45,45,45,45	0
34	NA	A	8334	1/1	0.95	0.07	30,30,30,30	0
32	MG	A	8057	1/1	0.95	0.09	31,31,31,31	0
32	MG	A	8117	1/1	0.95	0.07	40,40,40,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
34	NA	A	8370	1/1	0.95	0.24	58,58,58,58	0
32	MG	A	8044	1/1	0.95	0.15	51,51,51,51	0
34	NA	A	8317	1/1	0.95	0.04	29,29,29,29	0
34	NA	C	8345	1/1	0.95	0.06	52,52,52,52	0
35	CL	D	8519	1/1	0.95	0.12	57,57,57,57	0
32	MG	A	8108	1/1	0.95	0.16	79,79,79,79	0
34	NA	A	8344	1/1	0.95	0.07	29,29,29,29	0
32	MG	A	8086	1/1	0.95	0.07	47,47,47,47	0
32	MG	Z	8109	1/1	0.95	0.07	41,41,41,41	0
35	CL	P	8508	1/1	0.95	0.08	85,85,85,85	0
32	MG	A	8112	1/1	0.95	0.12	45,45,45,45	0
34	NA	A	8353	1/1	0.95	0.05	27,27,27,27	0
32	MG	A	8098	1/1	0.95	0.16	18,18,18,18	0
34	NA	M	8380	1/1	0.96	0.16	48,48,48,48	0
32	MG	A	8047	1/1	0.96	0.08	55,55,55,55	0
32	MG	L	8069	1/1	0.96	0.05	55,55,55,55	0
34	NA	A	8303	1/1	0.96	0.17	62,62,62,62	0
32	MG	A	8070	1/1	0.96	0.09	33,33,33,33	0
34	NA	U	8343	1/1	0.96	0.04	24,24,24,24	0
32	MG	A	8085	1/1	0.96	0.07	63,63,63,63	0
33	K	A	8200	1/1	0.96	0.10	66,66,66,66	0
34	NA	A	8342	1/1	0.96	0.07	28,28,28,28	0
35	CL	A	8513	1/1	0.96	0.06	44,44,44,44	0
35	CL	S	8506	1/1	0.96	0.05	44,44,44,44	0
32	MG	A	8100	1/1	0.96	0.11	72,72,72,72	0
35	CL	A	8515	1/1	0.96	0.17	68,68,68,68	0
32	MG	A	8097	1/1	0.97	0.11	30,30,30,30	0
32	MG	A	8027	1/1	0.97	0.04	45,45,45,45	0
34	NA	A	8320	1/1	0.97	0.08	31,31,31,31	0
35	CL	K	8501	1/1	0.97	0.04	58,58,58,58	0
35	CL	A	8516	1/1	0.97	0.06	45,45,45,45	0
35	CL	K	8521	1/1	0.97	0.05	44,44,44,44	0
32	MG	A	8022	1/1	0.97	0.05	53,53,53,53	0
36	CD	P	8405	1/1	0.97	0.04	82,82,82,82	0
32	MG	A	8072	1/1	0.98	0.05	60,60,60,60	0
32	MG	A	8008	1/1	0.98	0.05	37,37,37,37	0
32	MG	A	8048	1/1	0.98	0.04	39,39,39,39	0
32	MG	D	8055	1/1	0.98	0.04	34,34,34,34	0
32	MG	A	8059	1/1	0.98	0.04	32,32,32,32	0
34	NA	S	8338	1/1	0.98	0.04	38,38,38,38	0
32	MG	A	8062	1/1	0.98	0.05	64,64,64,64	0
32	MG	A	8023	1/1	0.98	0.04	33,33,33,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
32	MG	A	8110	1/1	0.98	0.05	20,20,20,20	0
32	MG	A	8001	1/1	0.98	0.04	16,16,16,16	0
32	MG	A	8016	1/1	0.98	0.04	46,46,46,46	0
32	MG	A	8113	1/1	0.98	0.08	25,25,25,25	0
32	MG	A	8068	1/1	0.98	0.03	57,57,57,57	0
32	MG	A	8115	1/1	0.98	0.04	41,41,41,41	0
32	MG	A	8040	1/1	0.98	0.10	49,49,49,49	0
34	NA	J	8309	1/1	0.98	0.03	18,18,18,18	0
36	CD	4	8404	1/1	0.98	0.02	62,62,62,62	0
34	NA	N	8347	1/1	0.99	0.06	12,12,12,12	0
32	MG	A	8083	1/1	0.99	0.04	35,35,35,35	0
32	MG	A	8084	1/1	0.99	0.06	49,49,49,49	0
32	MG	A	8009	1/1	0.99	0.06	37,37,37,37	0
32	MG	A	8005	1/1	0.99	0.03	49,49,49,49	0
32	MG	A	8012	1/1	0.99	0.05	29,29,29,29	0
32	MG	A	8028	1/1	0.99	0.03	41,41,41,41	0
32	MG	A	8030	1/1	0.99	0.03	19,19,19,19	0
32	MG	A	8032	1/1	0.99	0.07	30,30,30,30	0
32	MG	4	8078	1/1	0.99	0.04	45,45,45,45	0
32	MG	A	8056	1/1	0.99	0.05	37,37,37,37	0
32	MG	A	8033	1/1	0.99	0.03	27,27,27,27	0
32	MG	A	8035	1/1	0.99	0.06	38,38,38,38	0
32	MG	A	8060	1/1	0.99	0.03	29,29,29,29	0
32	MG	A	8061	1/1	0.99	0.04	22,22,22,22	0
32	MG	A	8037	1/1	0.99	0.05	34,34,34,34	0
32	MG	A	8063	1/1	0.99	0.08	78,78,78,78	0
32	MG	A	8038	1/1	0.99	0.03	16,16,16,16	0
32	MG	A	8039	1/1	0.99	0.04	47,47,47,47	0
32	MG	A	8013	1/1	0.99	0.02	39,39,39,39	0
32	MG	A	8014	1/1	0.99	0.05	18,18,18,18	0
32	MG	A	8007	1/1	0.99	0.05	14,14,14,14	0
32	MG	A	8071	1/1	0.99	0.03	75,75,75,75	0
32	MG	A	8043	1/1	0.99	0.04	23,23,23,23	0
32	MG	A	8018	1/1	0.99	0.03	49,49,49,49	0
32	MG	A	8019	1/1	0.99	0.06	27,27,27,27	0
32	MG	A	8077	1/1	0.99	0.06	20,20,20,20	0
32	MG	A	8079	1/1	0.99	0.04	39,39,39,39	0
32	MG	A	8080	1/1	0.99	0.04	55,55,55,55	0
32	MG	A	8020	1/1	0.99	0.03	38,38,38,38	0
36	CD	V	8401	1/1	0.99	0.03	59,59,59,59	0
36	CD	1	8403	1/1	0.99	0.02	60,60,60,60	0
36	CD	2	8402	1/1	0.99	0.03	59,59,59,59	0

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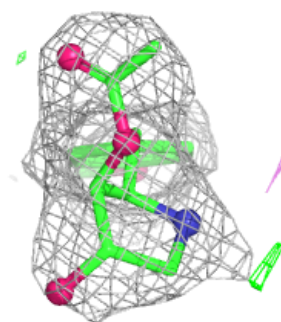
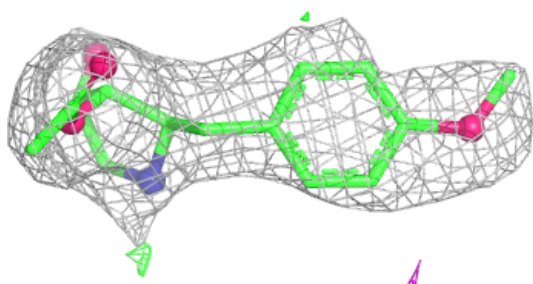
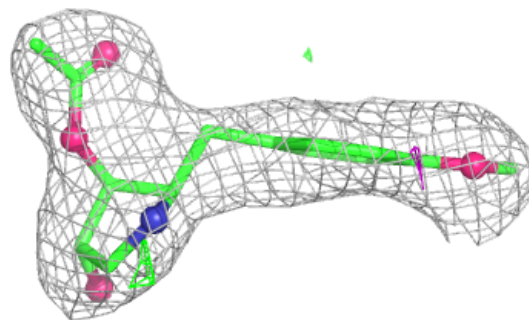
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
32	MG	A	8003	1/1	0.99	0.04	17,17,17,17	0
32	MG	A	8004	1/1	1.00	0.07	24,24,24,24	0
32	MG	A	8031	1/1	1.00	0.03	22,22,22,22	0
32	MG	A	8021	1/1	1.00	0.06	22,22,22,22	0
32	MG	A	8015	1/1	1.00	0.04	47,47,47,47	0
32	MG	C	8065	1/1	1.00	0.04	35,35,35,35	0
32	MG	A	8034	1/1	1.00	0.06	10,10,10,10	0
32	MG	A	8011	1/1	1.00	0.02	20,20,20,20	0
32	MG	A	8036	1/1	1.00	0.03	25,25,25,25	0
32	MG	A	8017	1/1	1.00	0.03	20,20,20,20	0
32	MG	A	8025	1/1	1.00	0.02	37,37,37,37	0
32	MG	A	8026	1/1	1.00	0.05	12,12,12,12	0
32	MG	A	8107	1/1	1.00	0.02	41,41,41,41	0
32	MG	A	8002	1/1	1.00	0.03	32,32,32,32	0
32	MG	A	8054	1/1	1.00	0.02	40,40,40,40	0
32	MG	A	8074	1/1	1.00	0.06	16,16,16,16	0
32	MG	A	8006	1/1	1.00	0.02	46,46,46,46	0
32	MG	A	8029	1/1	1.00	0.02	49,49,49,49	0
32	MG	A	8058	1/1	1.00	0.03	42,42,42,42	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ANM A 9000:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.