



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 11:01 PM GMT

PDB ID : 1VQ4
Title : The structure of the transition state analogue "DAA" bound to the large ribosomal subunit of *Haloarcula marismortui*
Authors : Schmeing, T.M.; Steitz, T.A.
Deposited on : 2004-12-16
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

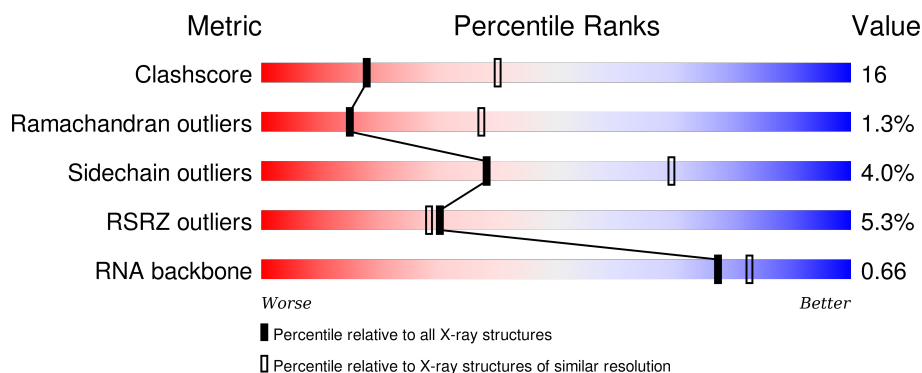
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.



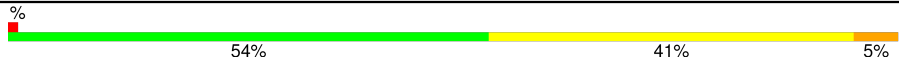
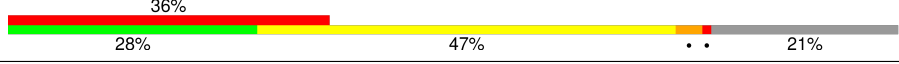
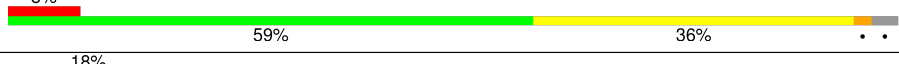

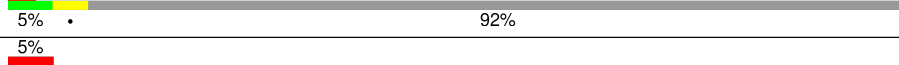
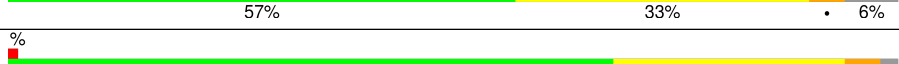
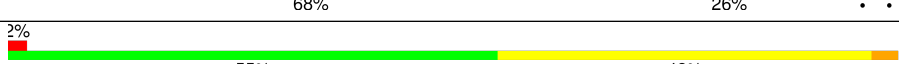

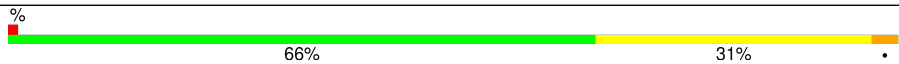


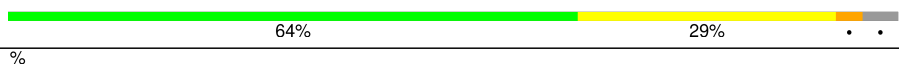
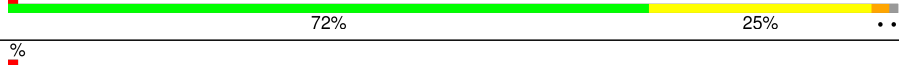


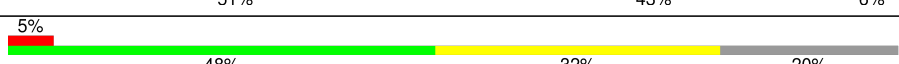


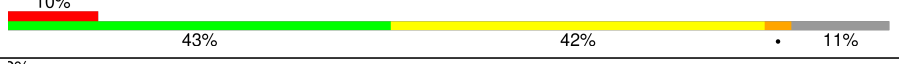
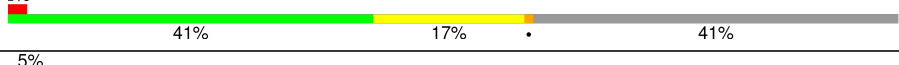





Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)
RNA backbone	2183	1069 (3.10-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	0	2922	<div> <div>2%</div> <div>59% 30% 5% • 6%</div> </div>
2	9	122	<div> <div>3%</div> <div>53% 34% 11% •</div> </div>
3	4	8	<div> <div>50% 38% 13%</div> </div>
4	A	240	<div> <div>4%</div> <div>56% 38% • •</div> </div>
5	B	338	<div> <div>2%</div> <div>52% 42% 6%</div> </div>

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Mol	Chain	Length	Quality of chain
6	C	246	
7	D	177	
8	E	178	
9	F	120	
10	G	348	
11	H	171	
12	J	145	
13	K	132	
14	L	165	
15	M	194	
16	N	187	
17	O	116	
18	P	149	
19	Q	96	
20	R	155	
21	S	85	
22	T	120	
23	U	66	
24	V	71	
25	W	154	
26	X	92	
27	Y	241	
28	Z	83	
29	1	57	
30	2	50	

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Mol	Chain	Length	Quality of chain
31	3	92	
32	I	162	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
33	MG	0	8053	-	-	-	X
33	MG	0	8060	-	-	-	X
34	K	0	9001	-	-	-	X
35	NA	0	9102	-	-	-	X
35	NA	0	9121	-	-	-	X
35	NA	0	9125	-	-	-	X
35	NA	0	9126	-	-	-	X
35	NA	0	9129	-	-	-	X
35	NA	0	9135	-	-	-	X
35	NA	0	9150	-	-	-	X
35	NA	0	9155	-	-	-	X
35	NA	0	9156	-	-	-	X
35	NA	0	9161	-	-	-	X
35	NA	0	9162	-	-	-	X
35	NA	0	9168	-	-	-	X
35	NA	0	9171	-	-	-	X
35	NA	0	9172	-	-	-	X
35	NA	0	9173	-	-	-	X
35	NA	0	9174	-	-	-	X
35	NA	0	9177	-	-	-	X
35	NA	0	9178	-	-	-	X
35	NA	0	9182	-	-	-	X
35	NA	L	9180	-	-	-	X
35	NA	M	9147	-	-	-	X
35	NA	Q	9148	-	-	-	X
35	NA	R	9186	-	-	-	X
36	CL	0	9315	-	-	-	X

2 Entry composition [i](#)

There are 38 unique types of molecules in this entry. The entry contains 98999 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 ribosomal rna.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	0	2754	Total	C	N	O	P	0	0	0
			59021	26350	10878	19048	2745			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
0	628	1MA	A	modified residue	GB 3377779
0	2587	OMU	U	modified residue	GB 3377779
0	2588	OMG	G	modified residue	GB 3377779
0	2619	UR3	U	modified residue	GB 3377779
0	2621	PSU	U	modified residue	GB 3377779

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	9	122	Total	C	N	O	P	0	0	0
			2600	1160	472	847	121			

- Molecule 3 is a RNA chain called 5'-R(*CP*CP*(5AA)P*(2OP)P*(PO2)P*(DA)P*C*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	4	8	Total	C	N	O	P	0	0	0
			127	61	23	38	5			

- Molecule 4 is a protein called 50S ribosomal protein L2P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	A	237	Total	C	N	O	S	0	0	0
			1753	1072	352	324	5			

- Molecule 5 is a protein called 50S ribosomal protein L3P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	B	337	Total	C	N	O	S	0	0	0
			2625	1616	493	511	5			

- Molecule 6 is a protein called 50S ribosomal protein L4E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	C	246	Total	C	N	O	S	0	0	0
			1859	1131	344	383	1			

- Molecule 7 is a protein called 50S ribosomal protein L5P.

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

- Molecule 8 is a protein called 50S ribosomal protein L6P.

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

- Molecule 9 is a protein called 50S ribosomal protein L7AE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	F	119	Total	C	N	O	S	0	0	0
			890	551	141	197	1			

- Molecule 10 is a protein called ACIDIC RIBOSOMAL PROTEIN P0 HOMOLOG.

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

- Molecule 11 is a protein called 50S RIBOSOMAL PROTEIN L10E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	H	160	Total	C	N	O	S	0	0	0
			1266	785	237	238	6			

- Molecule 12 is a protein called 50S ribosomal protein L13P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	J	142	Total	C	N	O	S	0	0	0
			1120	696	199	222	3			

- Molecule 13 is a protein called 50S ribosomal protein L14P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	K	132	Total	C	N	O	S	0	0	0
			992	609	187	192	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	44	LEU	HIS	CONFLICT	UNP P22450

- Molecule 14 is a protein called 50S ribosomal protein L15P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	L	145	Total	C	N	O	S	0	0	0
			1118	670	222	226				

- Molecule 15 is a protein called 50S Ribosomal Protein L15E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	M	194	Total	C	N	O	S	0	0	0
			1560	943	332	284	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	13	GLU	LYS	CONFLICT	GB 55231501
M	194	ALA	GLY	CONFLICT	GB 55231501

- Molecule 16 is a protein called 50S ribosomal protein L18P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	N	186	Total	C	N	O	S	0	0	0
			1445	895	262	286	2			

- Molecule 17 is a protein called 50S ribosomal protein L18e.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
17	O	115	Total	C	N	O	0	0	0
			865	529	161	175			

- Molecule 18 is a protein called 50S ribosomal protein L19E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	P	143	Total	C	N	O	0	0	0
			1136	683	229	224			

- Molecule 19 is a protein called 50S ribosomal protein L21e.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
19	Q	95	Total	C	N	O	0	0	0
			735	450	141	144			

- Molecule 20 is a protein called 50S ribosomal protein L22P.

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

- Molecule 21 is a protein called 50S ribosomal protein L23P.

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

- Molecule 22 is a protein called 50S ribosomal protein L24P.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
22	T	119	Total	C	N	O	0	0	0
			950	568	180	202			

- Molecule 23 is a protein called 50S ribosomal protein L24E.

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

- Molecule 24 is a protein called 50S ribosomal protein L29P.

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

- Molecule 25 is a protein called 50S ribosomal protein L30P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	W	154	Total	C	N	O	S	0	0	0
			1196	737	209	244	6			

- Molecule 26 is a protein called 50S ribosomal protein L31e.

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

- Molecule 27 is a protein called 50S ribosomal protein L32E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	Y	142	Total	C	N	O	S	0	0	0
			1130	686	228	216				

- Molecule 28 is a protein called 50S ribosomal protein L37Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	Z	73	Total	C	N	O	S	0	0	0
			578	346	116	111	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Z	10	ARG	SER	CONFLICT	GB 55231162

- Molecule 29 is a protein called 50S ribosomal protein L37e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	1	56	Total	C	N	O	S	0	0	0
			431	258	86	83	4			

- Molecule 30 is a protein called 50S ribosomal protein L39e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	2	46	Total	C	N	O	S	0	0	0
			396	239	89	67	1			

- Molecule 31 is a protein called 50S ribosomal protein L44E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	3	92	Total	C	N	O	S	0	0	0
			755	458	153	137	7			

- Molecule 32 is a protein called 50S RIBOSOMAL PROTEIN L11P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	I	70	Total	C	N	O	S	0	0	0
			519	323	81	114	1			

- Molecule 33 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
33	0	108	Total	Mg	0	0
			108	108		
33	Y	1	Total	Mg	0	0
			1	1		
33	K	1	Total	Mg	0	0
			1	1		
33	B	1	Total	Mg	0	0
			1	1		
33	A	2	Total	Mg	0	0
			2	2		
33	T	1	Total	Mg	0	0
			1	1		
33	9	2	Total	Mg	0	0
			2	2		
33	3	1	Total	Mg	0	0
			1	1		

- Molecule 34 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
34	0	3	Total	K	0	0
			3	3		

- Molecule 35 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
35	0	72	Total Na 72 72	0	0
35	J	1	Total Na 1 1	0	0
35	Q	1	Total Na 1 1	0	0
35	H	2	Total Na 2 2	0	0
35	C	1	Total Na 1 1	0	0
35	A	1	Total Na 1 1	0	0
35	R	3	Total Na 3 3	0	0
35	9	2	Total Na 2 2	0	0
35	L	1	Total Na 1 1	0	0
35	S	1	Total Na 1 1	0	0
35	M	1	Total Na 1 1	0	0

- Molecule 36 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
36	0	10	Total Cl 10 10	0	0
36	J	3	Total Cl 3 3	0	0
36	K	1	Total Cl 1 1	0	0
36	B	1	Total Cl 1 1	0	0
36	A	1	Total Cl 1 1	0	0
36	N	1	Total Cl 1 1	0	0
36	O	1	Total Cl 1 1	0	0
36	R	1	Total Cl 1 1	0	0
36	L	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
36	3	1	Total 1	Cl 1	0	0
36	M	1	Total 1	Cl 1	0	0

- Molecule 37 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	O	1	Total 1	Cd 1	0	0
37	Z	1	Total 1	Cd 1	0	0
37	1	1	Total 1	Cd 1	0	0
37	3	1	Total 1	Cd 1	0	0
37	U	1	Total 1	Cd 1	0	0

- Molecule 38 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
38	0	5764	Total 5764	O 5764	0	0
38	9	133	Total 133	O 133	0	0
38	4	3	Total 3	O 3	0	0
38	A	116	Total 116	O 116	0	0
38	B	143	Total 143	O 143	0	0
38	C	173	Total 173	O 173	0	0
38	D	44	Total 44	O 44	0	0
38	E	43	Total 43	O 43	0	0
38	F	24	Total 24	O 24	0	0
38	G	17	Total 17	O 17	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
38	H	66	Total 66	O 66	0	0
38	J	52	Total 52	O 52	0	0
38	K	57	Total 57	O 57	0	0
38	L	81	Total 81	O 81	0	0
38	M	115	Total 115	O 115	0	0
38	N	61	Total 61	O 61	0	0
38	O	45	Total 45	O 45	0	0
38	P	63	Total 63	O 63	0	0
38	Q	52	Total 52	O 52	0	0
38	R	89	Total 89	O 89	0	0
38	S	31	Total 31	O 31	0	0
38	T	36	Total 36	O 36	0	0
38	U	26	Total 26	O 26	0	0
38	V	13	Total 13	O 13	0	0
38	W	70	Total 70	O 70	0	0
38	X	31	Total 31	O 31	0	0
38	Y	93	Total 93	O 93	0	0
38	Z	31	Total 31	O 31	0	0
38	1	61	Total 61	O 61	0	0
38	2	42	Total 42	O 42	0	0
38	3	71	Total 71	O 71	0	0

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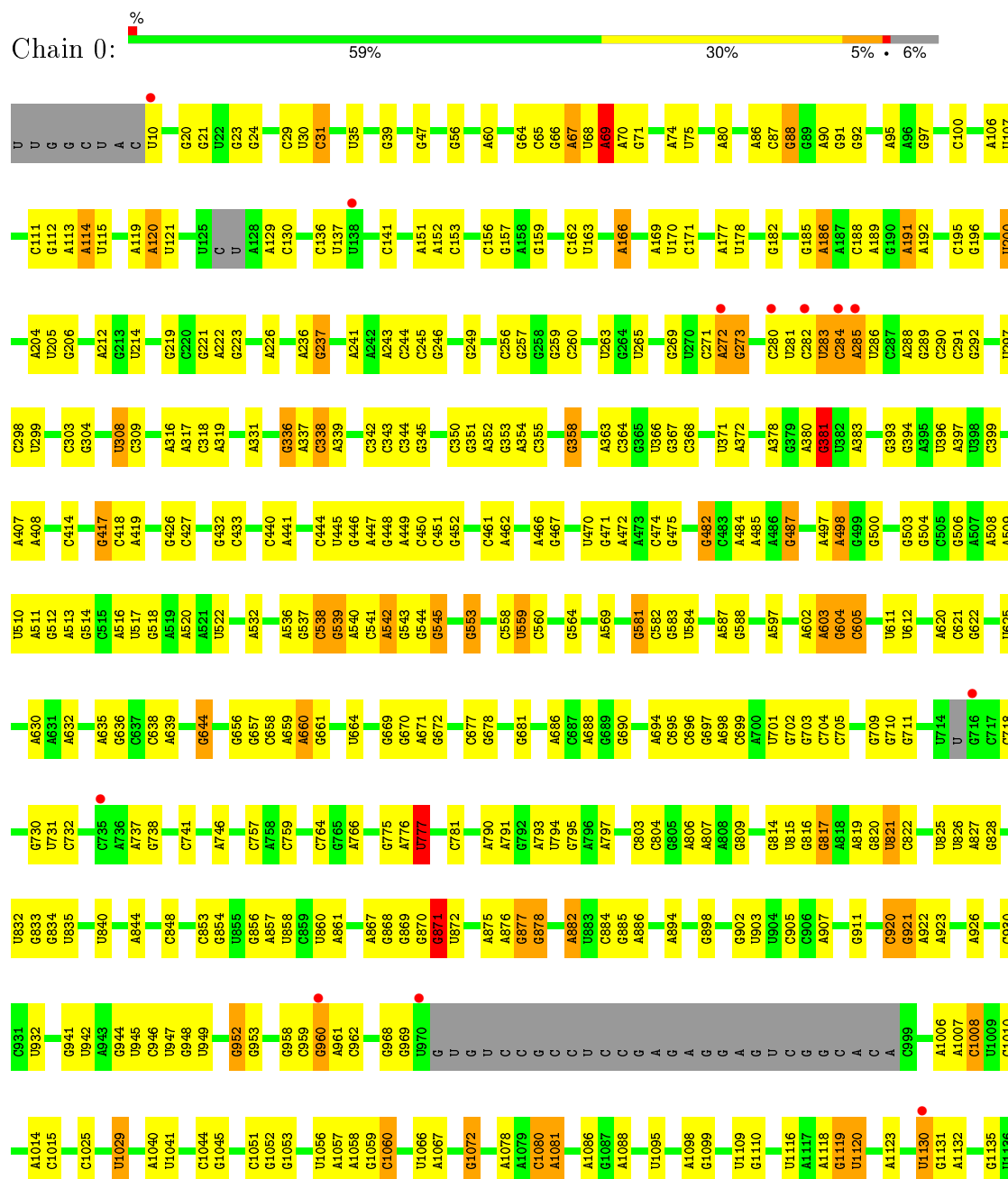
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
38	I	9	Total	O	0	0
			9	9		

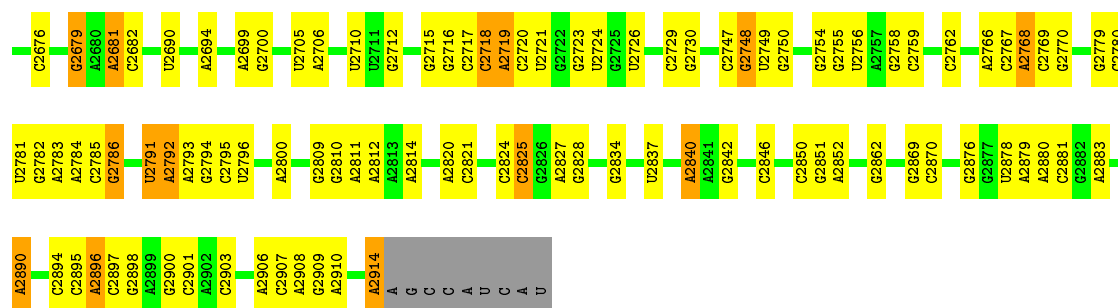
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

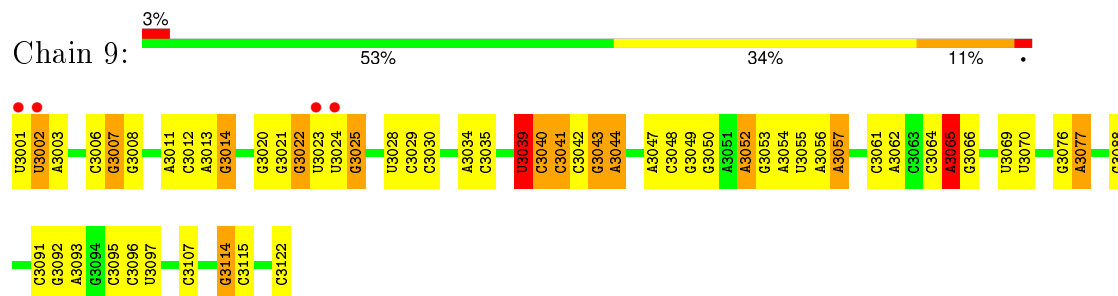
- Molecule 1: 23S ribosomal rna



C2561	C2562	C2563	C2564	C2565	C2570	C2578	C2586	C2587	C2588	C2589	C2590	C2591	C2592	C2597	C2598	A2601	A2602	A2603	A2604	A2605	A2606	C2607	C2608	C2613	C2614	C2619	C2626	C2627	C2630	C2634	C2635	C2636	C2637	C2638	C2642	C2643	C2644	C2649	C2652	A2664	A	U	C2667	C2668	C2669	C2670	C2671	C2672	U2673
A2469	C2472	U2473	C2474	C2475	C2476	C2477	C2478	A2479	C2480	A2483	C2489	A2490	C2491	U2492	C2493	C2502	A2503	A2504	C2505	A2506	C2507	C2508	A2509	C2510	A2511	C2515	C2516	A2521	C2524	C2525	C2526	C2527	U2531	A2532	C2533	C2534	U2535	C2536	C2537	U2541	C2542	U2545	C2546	C2547	C2548	C2552	C2553	U2673	
A2353	A2354	A2361	A2362	C2363	C2364	C2365	A2369	A2379	A2380	C2381	A2382	A2385	C2403	C2404	A2408	C2412	A2413	A2414	A2415	C2416	C2417	C2418	U2419	C2420	C2421	U2422	C2423	U2424	A2425	C2426	C2427	C2428	A2434	U2435	U2436	U2441	C2442	C2443	G2451	C2452	C2453	C2462	A2465	C2466	A2468				
C2241	U2242	C2243	C2249	C2250	C2251	C2252	C2253	C2256	C2257	A2258	A2264	U2265	A2266	C2269	C2270	C2271	C2272	C2281	C2282	A2291	A2300	A2301	A2302	C2309	C2312	C2313	C2314	C2315	C2316	C2317	U2320	A2321	U2322	C2323	C2324	C2325	U2326	A2332	C2338	A	C	A	C	A	C2344	A2346			
U1984	A1989	G1970	G1971	U1972	A1973	A1978	C1979	U1980	U1986	U2003	U2004	U2008	C2009	A2010	A2011	U2012	C2013	C2014	A2015	U2016	U2008	C2009	A2010	A2011	U2012	C2013	C2014	A2015	U2016	U2008	C2009	A2010	A2011	U2012	C2013	C2014	A2015	U2016	U2008	C2009	A2010	A2011	U2012	C2013	C2014	A2015	U2016		
C1862	G1863	G1867	G1868	U1874	C1877	G1877	U1878	U1879	U1890	U1903	A1904	U1905	A1909	A1910	A1921	G1925	G1926	A1927	U1937	U1938	U1939	C1940	A1941	C1942	G1943	G1944	C1946	G1947	G1948	G1949	G1950	G1951	U	A	C	C	U	A	C	U	A	C	U	A	C				
G1666	A1667	U1668	A1669	G1670	C1675	G1676	U1677	G1681	A1682	G1683	A1684	A1685	C1686	C1687	C1692	A1701	U1702	C1705	A1710	A1711	A1712	A1717	U1722	G1723	C1725	G1730	C1731	A1732	A1733	C1734	C1735	U1741	A1742	G1743	G1744	A1747	G1751	G1752	A1755	A1756	A1759	C1764							
G1557	C1558	A1559	U	U1561	C1562	C1563	C1564	U1568	U1569	A1572	A1573	C1574	G1589	G1592	C1593	C1594	U1595	A1597	A1598	A1603	G1604	G1605	A1606	A1607	C1613	G1614	C1617	U1625	A1626	A1630	A1631	A1632	C1633	G1634	U1635	G1636	A1637	A1641	A1642	C1643	U1654	A1655	A1656	A1657	A1658	G1660			
C1439	U1440	G1441	A1442	C1450	C1451	C1462	A1463	A1471	C1472	U1473	C1474	C1477	U1478	G1481	A1482	A1485	A1486	A1494	C1495	G1496	U1500	U1503	A1504	U1505	U1506	C1507	U1517	C1521	A1522	G1523	U1524	G1525	A1526	A1527	A1528	G1529	G1535	C1536	C1537	G1543	U1544	C1545	G1546	G1555	G1556				
A1329	U1333	C1334	C1335	G1340	A1341	C1342	C1343	G1344	G1351	C1352	C1353	C1360	C1366	A1372	A1375	G1376	C1377	G1385	U1388	G1389	A1390	G1391	A1392	A1393	C1394	G1398	A1399	A1406	A1407	A1414	G1415	G1416	G1417	U1418	U1419	U1422	C1423	A1424	G1425	C1426	A1427	A1434	U1435	C1436					
C1229	A1230	U1234	C1235	A1236	U1237	C1238	G1239	A1242	C1243	U1244	C1245	A1246	A1252	C1253	U1266	C1267	C1268	C1273	A1278	U1279	U1285	C1289	G1290	U1293	A1294	G1295	U1298	G1299	U1304	C1305	U1306	A1307	A1308	G1311	A1312	A1313	U1314	G1315	G1316	A1321	G1322	G1325	A1328						
G1137	G1151	G1156	C1157	G1158	G1159	G1160	A1161	G1162	G1163	U1164	G1165	A1166	G1167	C1168	U1169	U1170	C1172	A1173	G1175	C1176	A1177	U1180	A1181	C1182	C1183	U1185	U1187	A1188	A1189	G1190	A1191	A1192	A1193	U1198	A1199	A1200	C1201	G1205	U1206	A1207	C1208	C1209	G1210	C1213	G1214	G1216			



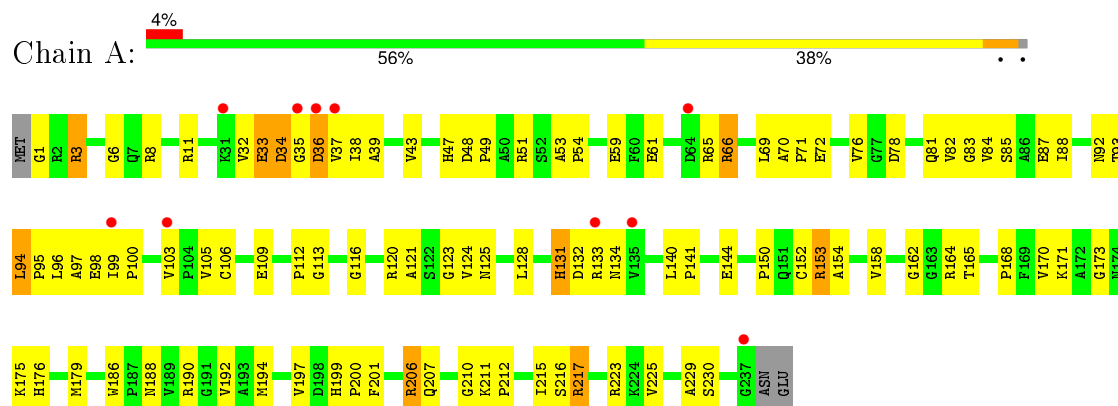
• Molecule 2: 5S ribosomal RNA



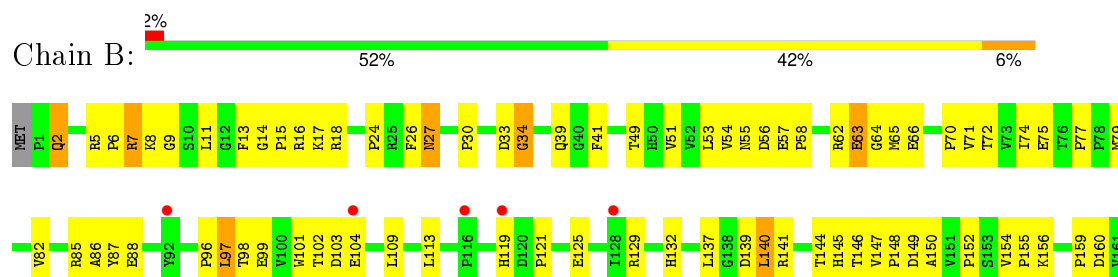
• Molecule 3: 5'-R(*CP*CP*(5AA)P*(2OP)P*(PO2)P*(DA)P*C*C)-3''

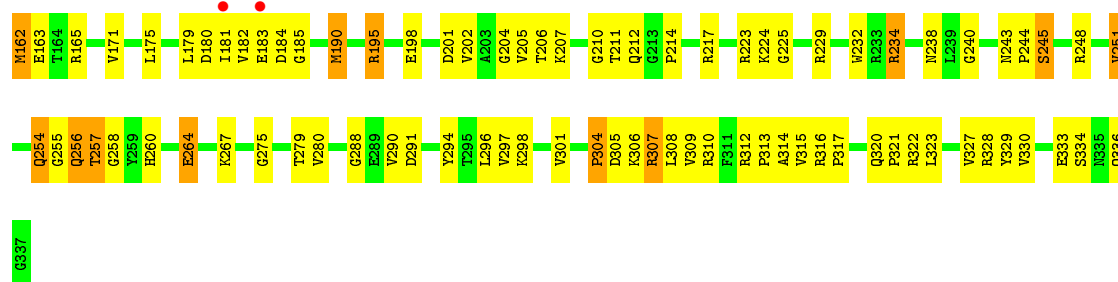


• Molecule 4: 50S ribosomal protein L2P

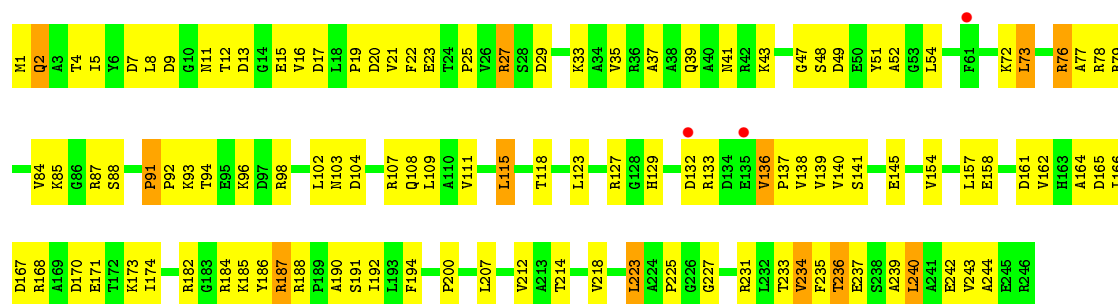


• Molecule 5: 50S ribosomal protein L3P

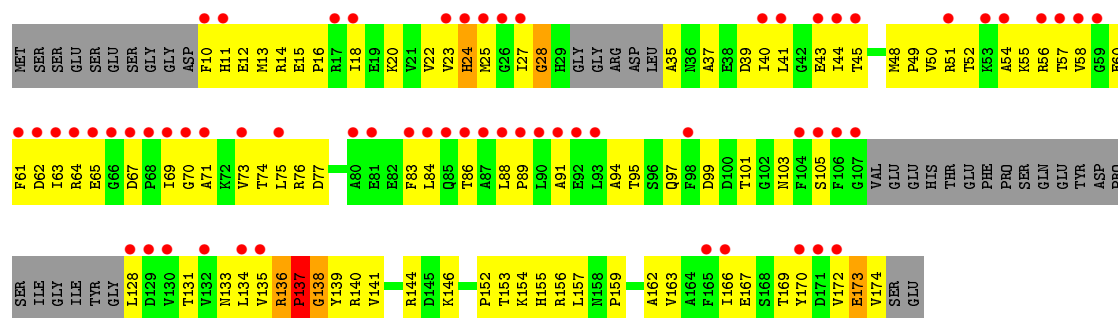




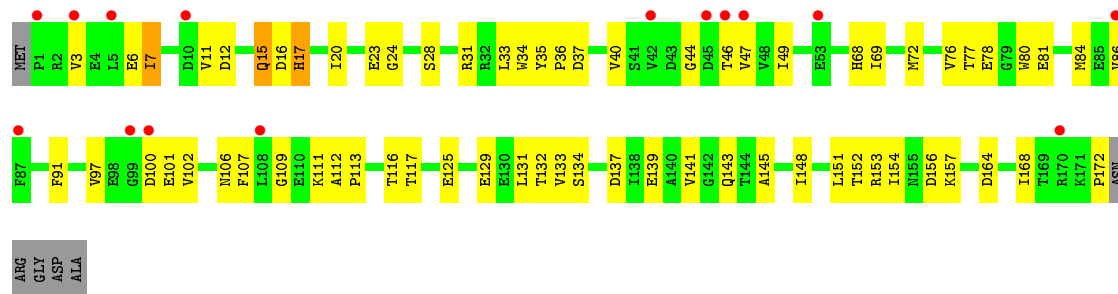
- Molecule 6: 50S ribosomal protein L4E



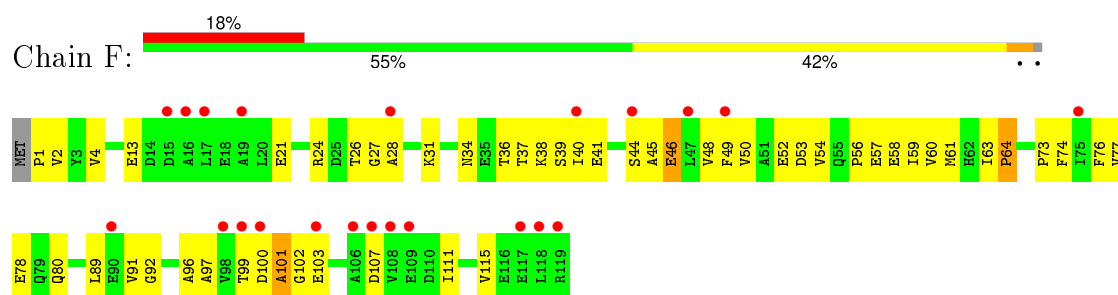
- Molecule 7: 50S ribosomal protein L5P



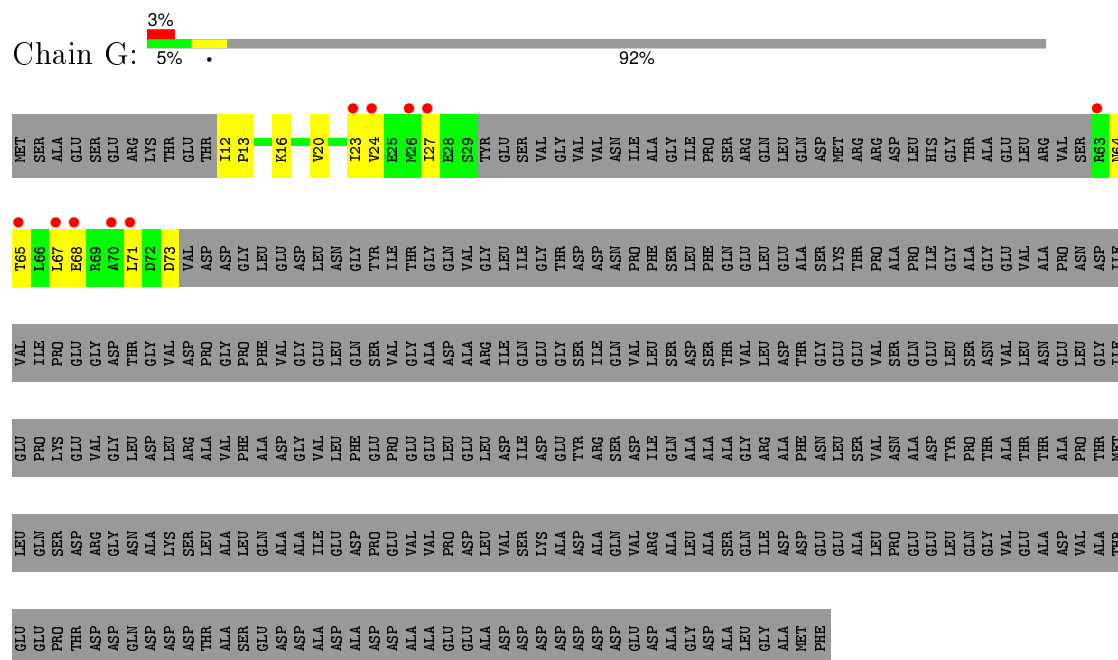
- Molecule 8: 50S ribosomal protein L6P



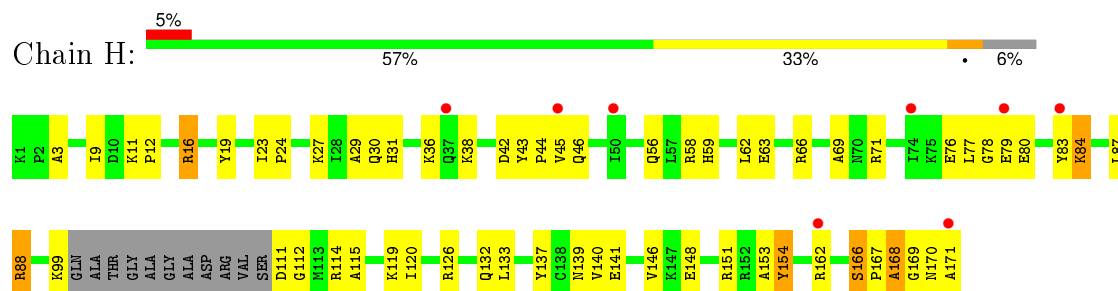
- Molecule 9: 50S ribosomal protein L7AE



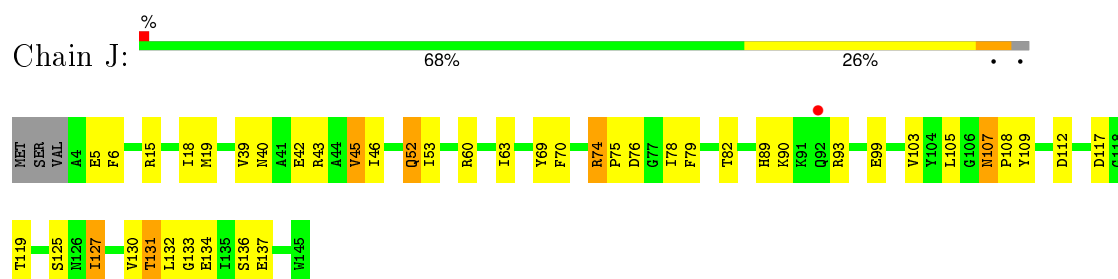
• Molecule 10: ACIDIC RIBOSOMAL PROTEIN P0 HOMOLOG



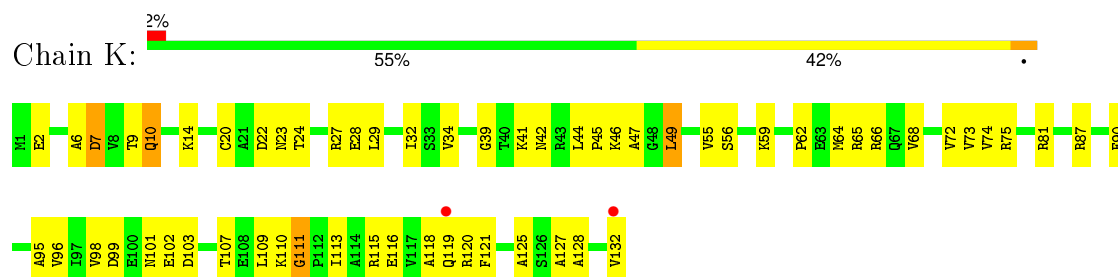
- Molecule 11: 50S RIBOSOMAL PROTEIN L10E



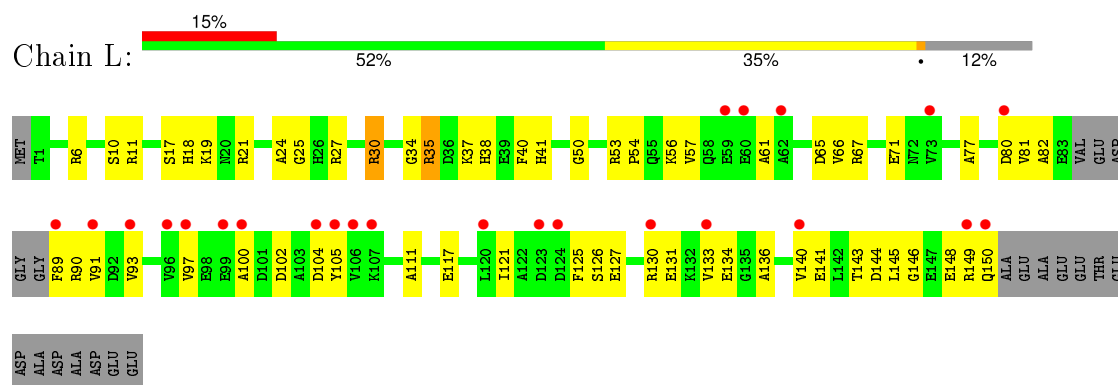
- Molecule 12: 50S ribosomal protein L13P



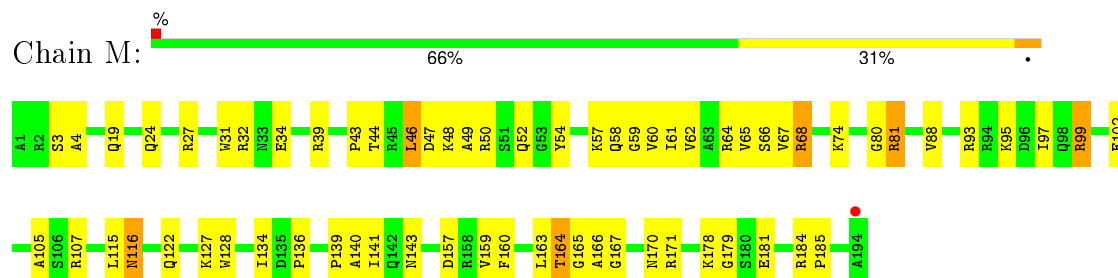
- Molecule 13: 50S ribosomal protein L14P



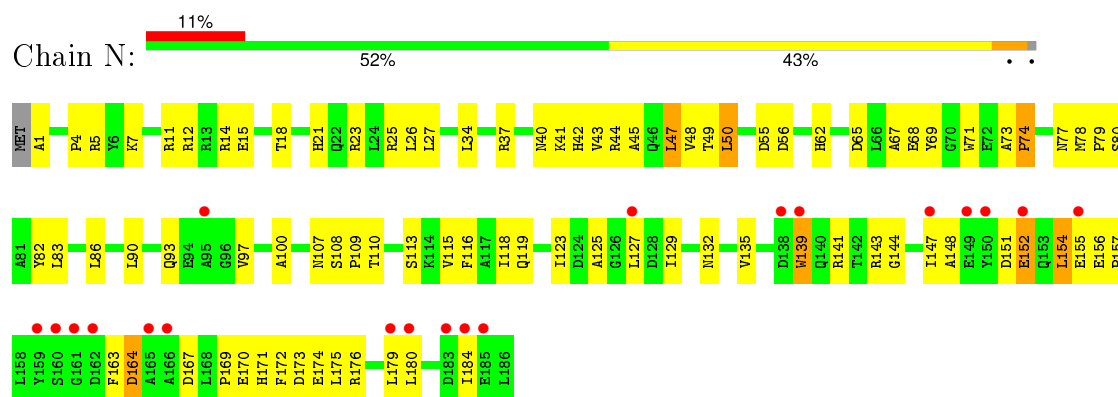
- Molecule 14: 50S ribosomal protein L15P



- Molecule 15: 50S Ribosomal Protein L15E



- Molecule 16: 50S ribosomal protein L18P



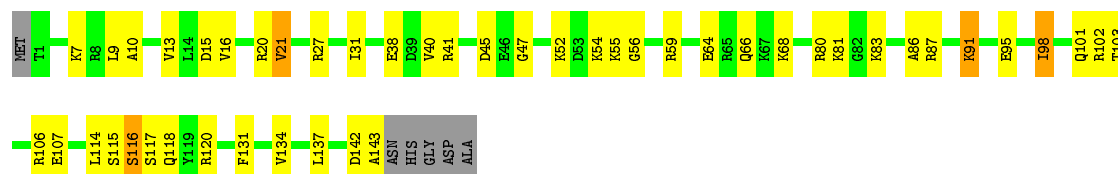
- Molecule 17: 50S ribosomal protein L18e





- Molecule 18: 50S ribosomal protein L19E

Chain P: 64% 29%



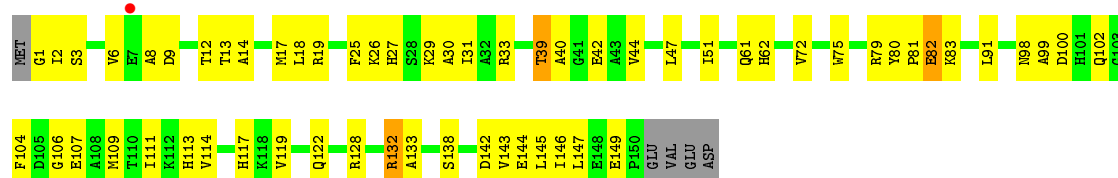
- Molecule 19: 50S ribosomal protein L21e

Chain Q: 72% 25%



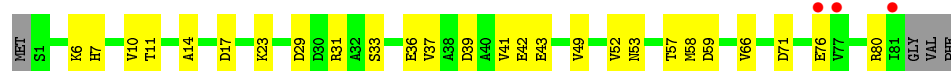
- Molecule 20: 50S ribosomal protein L22P

Chain R: 58% 37%



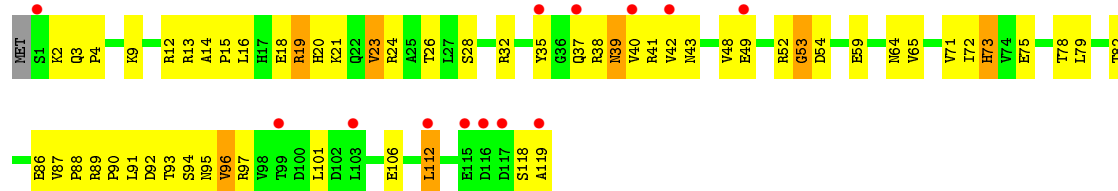
- Molecule 21: 50S ribosomal protein L23P

Chain S: 65% 31% 5%



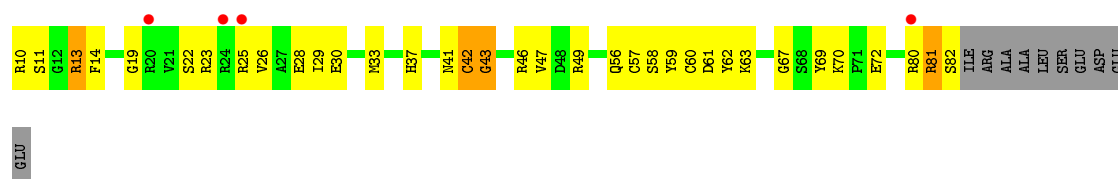
- Molecule 22: 50S ribosomal protein L24P

Chain T: 51% 43% 6%



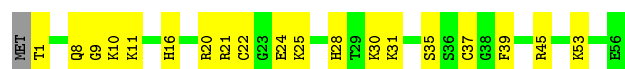
- Molecule 23: 50S ribosomal protein L24E





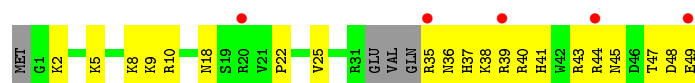
- Molecule 29: 50S ribosomal protein L37e

Chain 1: 65% 33%



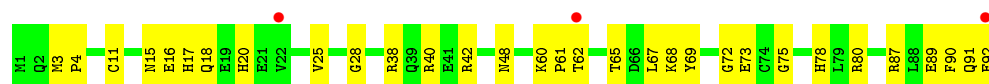
- Molecule 30: 50S ribosomal protein L39e

Chain 2: 10% 50% 42% 8%



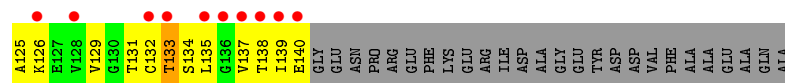
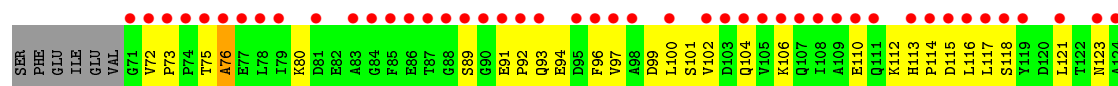
- Molecule 31: 50S ribosomal protein L44E

Chain 3: 3% 66% 34%



- Molecule 32: 50S RIBOSOMAL PROTEIN L11P

Chain I: 19% 35% 23% 57%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	213.00Å 301.03Å 575.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 49.83 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.70) 94.8 (49.83-2.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.69Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.190 , 0.230 0.183 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	54.6	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 505940 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	98999	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.58% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, OMG, CL, NA, K, PO2, CD, 5AA, OMU, UR3, 2OP, 1MA, PSU

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	0	0.38	0/65959	0.69	20/102870 (0.0%)
2	9	0.35	0/2905	0.71	2/4528 (0.0%)
3	4	0.46	0/102	0.73	0/149
4	A	0.31	0/1786	0.65	0/2408
5	B	0.34	0/2690	0.65	0/3652
6	C	0.37	0/1884	0.65	1/2551 (0.0%)
7	D	0.30	0/1111	0.53	0/1498
8	E	0.33	0/1382	0.58	0/1880
9	F	0.30	0/901	0.54	0/1224
10	G	0.30	0/241	0.49	0/324
11	H	0.35	0/1287	0.67	0/1725
12	J	0.37	0/1136	0.63	0/1530
13	K	0.36	0/1001	0.69	0/1347
14	L	0.33	0/1130	0.64	0/1509
15	M	0.33	0/1584	0.61	0/2119
16	N	0.30	0/1474	0.63	0/1999
17	O	0.31	0/874	0.56	0/1181
18	P	0.33	0/1147	0.54	0/1528
19	Q	0.36	0/749	0.72	0/1005
20	R	0.36	0/1172	0.65	0/1578
21	S	0.32	0/648	0.57	0/875
22	T	0.32	0/958	0.61	0/1289
23	U	0.34	0/417	0.58	0/562
24	V	0.27	0/502	0.54	0/675
25	W	0.36	0/1219	0.63	0/1655
26	X	0.36	0/664	0.58	0/895
27	Y	0.37	0/1146	0.64	0/1536
28	Z	0.36	0/589	0.67	0/787
29	1	0.36	0/438	0.62	0/578
30	2	0.33	0/401	0.52	0/529
31	3	0.38	0/771	0.58	0/1024
32	I	0.30	0/526	0.54	0/716

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.37	0/98794	0.67	23/147726 (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	0	0	47
2	9	0	2
3	4	0	1
25	W	0	1
All	All	0	51

There are no bond length outliers.

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0	1563	G	C2'-C3'-O3'	8.54	128.28	109.50
1	0	1942	A	C5'-C4'-C3'	8.16	129.06	116.00
1	0	1979	G	C2'-C3'-O3'	6.72	124.45	113.70
1	0	871	G	C5'-C4'-O4'	-6.61	101.17	109.10
2	9	3039	U	N1-C1'-C2'	6.29	122.18	114.00

There are no chirality outliers.

5 of 51 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	0	191	A	Sidechain
1	0	221	G	Sidechain
1	0	24	G	Sidechain
1	0	417	G	Sidechain
1	0	471	G	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	0	59021	0	29812	849	0
2	9	2600	0	1326	65	0
3	4	127	0	75	4	0
4	A	1753	0	1766	119	0
5	B	2625	0	2533	159	0
6	C	1859	0	1816	127	0
7	D	1094	0	1085	91	0
8	E	1357	0	1266	64	0
9	F	890	0	843	56	0
10	G	240	0	231	13	0
11	H	1266	0	1268	70	0
12	J	1120	0	1098	55	0
13	K	992	0	1031	65	0
14	L	1118	0	1076	55	0
15	M	1560	0	1568	70	0
16	N	1445	0	1401	107	0
17	O	865	0	873	39	0
18	P	1136	0	1123	44	0
19	Q	735	0	729	23	0
20	R	1149	0	1122	62	0
21	S	641	0	605	21	0
22	T	950	0	923	61	0
23	U	410	0	364	24	0
24	V	499	0	511	33	0
25	W	1196	0	1137	95	0
26	X	654	0	653	50	0
27	Y	1130	0	1133	55	0
28	Z	578	0	539	27	0
29	1	431	0	426	30	0
30	2	396	0	413	27	0
31	3	755	0	728	31	0
32	I	519	0	500	54	0
33	0	108	0	0	0	0
33	3	1	0	0	0	0
33	9	2	0	0	0	0
33	A	2	0	0	0	0
33	B	1	0	0	0	0
33	K	1	0	0	0	0
33	T	1	0	0	0	0
33	Y	1	0	0	0	0
34	0	3	0	0	0	0
35	0	72	0	0	0	0
35	9	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
35	A	1	0	0	0	0
35	C	1	0	0	0	0
35	H	2	0	0	0	0
35	J	1	0	0	0	0
35	L	1	0	0	0	0
35	M	1	0	0	0	0
35	Q	1	0	0	0	0
35	R	3	0	0	0	0
35	S	1	0	0	0	0
36	0	10	0	0	1	0
36	3	1	0	0	0	0
36	A	1	0	0	0	0
36	B	1	0	0	0	0
36	J	3	0	0	1	0
36	K	1	0	0	0	0
36	L	1	0	0	0	0
36	M	1	0	0	1	0
36	N	1	0	0	1	0
36	O	1	0	0	0	0
36	R	1	0	0	0	0
37	1	1	0	0	0	0
37	3	1	0	0	0	0
37	O	1	0	0	0	0
37	U	1	0	0	0	0
37	Z	1	0	0	0	0
38	0	5764	0	0	97	0
38	1	61	0	0	3	0
38	2	42	0	0	3	0
38	3	71	0	0	5	0
38	4	3	0	0	0	0
38	9	133	0	0	4	0
38	A	116	0	0	18	0
38	B	143	0	0	23	0
38	C	173	0	0	21	0
38	D	44	0	0	8	0
38	E	43	0	0	5	0
38	F	24	0	0	4	0
38	G	17	0	0	0	0
38	H	66	0	0	9	0
38	I	9	0	0	2	0
38	J	52	0	0	3	0
38	K	57	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
38	L	81	0	0	11	0
38	M	115	0	0	4	0
38	N	61	0	0	10	0
38	O	45	0	0	6	0
38	P	63	0	0	3	0
38	Q	52	0	0	1	0
38	R	89	0	0	5	0
38	S	31	0	0	2	0
38	T	36	0	0	2	0
38	U	26	0	0	0	0
38	V	13	0	0	1	0
38	W	70	0	0	5	0
38	X	31	0	0	5	0
38	Y	93	0	0	7	0
38	Z	31	0	0	1	0
All	All	98999	0	59974	2378	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 16.

The worst 5 of 2378 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)
2:9:3006:C:H5''	16:N:37:ARG:NH1	1.64	1.13
6:C:236:THR:HG22	6:C:239:ALA:H	1.11	1.13
11:H:46:GLN:HB3	11:H:167:PRO:HD2	1.32	1.11
2:9:3006:C:H5''	16:N:37:ARG:HH12	1.08	1.07
1:0:1160:G:H5'	1:0:1161:A:H5'	1.34	1.04

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
4	A	235/240 (98%)	205 (87%)	28 (12%)	2 (1%)	21	49
5	B	335/338 (99%)	303 (90%)	27 (8%)	5 (2%)	13	32
6	C	244/246 (99%)	223 (91%)	20 (8%)	1 (0%)	39	69
7	D	134/177 (76%)	94 (70%)	34 (25%)	6 (4%)	3	6
8	E	170/178 (96%)	161 (95%)	7 (4%)	2 (1%)	16	39
9	F	117/120 (98%)	101 (86%)	13 (11%)	3 (3%)	7	16
10	G	25/348 (7%)	24 (96%)	1 (4%)	0	100	100
11	H	156/171 (91%)	144 (92%)	7 (4%)	5 (3%)	5	12
12	J	140/145 (97%)	128 (91%)	10 (7%)	2 (1%)	14	35
13	K	130/132 (98%)	116 (89%)	12 (9%)	2 (2%)	13	32
14	L	141/165 (86%)	118 (84%)	23 (16%)	0	100	100
15	M	192/194 (99%)	178 (93%)	13 (7%)	1 (0%)	34	63
16	N	184/187 (98%)	167 (91%)	12 (6%)	5 (3%)	6	16
17	O	113/116 (97%)	107 (95%)	6 (5%)	0	100	100
18	P	141/149 (95%)	131 (93%)	9 (6%)	1 (1%)	26	55
19	Q	93/96 (97%)	90 (97%)	3 (3%)	0	100	100
20	R	148/155 (96%)	133 (90%)	14 (10%)	1 (1%)	26	55
21	S	79/85 (93%)	75 (95%)	4 (5%)	0	100	100
22	T	117/120 (98%)	108 (92%)	8 (7%)	1 (1%)	21	49
23	U	51/66 (77%)	47 (92%)	3 (6%)	1 (2%)	9	24
24	V	63/71 (89%)	55 (87%)	7 (11%)	1 (2%)	12	30
25	W	152/154 (99%)	144 (95%)	6 (4%)	2 (1%)	15	37
26	X	80/92 (87%)	72 (90%)	7 (9%)	1 (1%)	15	37
27	Y	140/241 (58%)	134 (96%)	6 (4%)	0	100	100
28	Z	71/83 (86%)	57 (80%)	9 (13%)	5 (7%)	1	2
29	1	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
30	2	42/50 (84%)	40 (95%)	2 (5%)	0	100	100
31	3	90/92 (98%)	85 (94%)	5 (6%)	0	100	100
32	I	68/162 (42%)	55 (81%)	11 (16%)	2 (3%)	6	14
All	All	3705/4430 (84%)	3348 (90%)	308 (8%)	49 (1%)	15	37

5 of 49 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	D	27	ILE
7	D	137	PRO
7	D	173	GLU
9	F	101	ALA
11	H	140	VAL

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	
4	A	179/182 (98%)	167 (93%)	12 (7%)	20	44
5	B	282/283 (100%)	263 (93%)	19 (7%)	20	44
6	C	193/193 (100%)	180 (93%)	13 (7%)	20	44
7	D	117/148 (79%)	112 (96%)	5 (4%)	35	66
8	E	152/156 (97%)	148 (97%)	4 (3%)	54	83
9	F	93/94 (99%)	92 (99%)	1 (1%)	80	94
10	G	27/283 (10%)	27 (100%)	0	100	100
11	H	132/138 (96%)	126 (96%)	6 (4%)	34	65
12	J	118/121 (98%)	109 (92%)	9 (8%)	16	37
13	K	106/106 (100%)	103 (97%)	3 (3%)	51	81
14	L	113/127 (89%)	108 (96%)	5 (4%)	35	65
15	M	158/158 (100%)	151 (96%)	7 (4%)	35	65
16	N	149/150 (99%)	145 (97%)	4 (3%)	52	82
17	O	93/94 (99%)	92 (99%)	1 (1%)	80	94
18	P	113/117 (97%)	109 (96%)	4 (4%)	43	74
19	Q	79/80 (99%)	76 (96%)	3 (4%)	40	71
20	R	117/122 (96%)	114 (97%)	3 (3%)	54	83
21	S	71/74 (96%)	69 (97%)	2 (3%)	51	81
22	T	105/106 (99%)	98 (93%)	7 (7%)	20	44
23	U	44/52 (85%)	44 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
24	V	51/57 (90%)	51 (100%)	0	100	100
25	W	130/130 (100%)	123 (95%)	7 (5%)	27	56
26	X	66/74 (89%)	62 (94%)	4 (6%)	23	49
27	Y	120/196 (61%)	116 (97%)	4 (3%)	45	76
28	Z	60/68 (88%)	59 (98%)	1 (2%)	68	90
29	1	46/47 (98%)	46 (100%)	0	100	100
30	2	42/46 (91%)	41 (98%)	1 (2%)	57	85
31	3	79/79 (100%)	79 (100%)	0	100	100
32	I	58/130 (45%)	58 (100%)	0	100	100
All	All	3093/3611 (86%)	2968 (96%)	125 (4%)	38	69

5 of 125 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
11	H	88	ARG
13	K	7	ASP
26	X	15	ARG
11	H	132	GLN
12	J	74	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 86 such sidechains are listed below:

Mol	Chain	Res	Type
15	M	58	GLN
19	Q	40	HIS
30	2	16	ASN
15	M	143	ASN
16	N	107	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	0	2745/2922 (93%)	237 (8%)	34 (1%)
2	9	121/122 (99%)	17 (14%)	1 (0%)
3	4	1/8 (12%)	0	0
All	All	2867/3052 (93%)	254 (8%)	35 (1%)

5 of 254 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	0	31	C
1	0	60	A
1	0	67	A
1	0	69	A
1	0	70	A

5 of 35 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	0	1246	A
1	0	1506	U
1	0	2718	C
1	0	1352	A
1	0	1377	C

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	OMU	0	2587	1	12,22,23	0.98	1 (8%)	19,31,34	3.19	2 (10%)
1	OMG	0	2588	1,3	17,26,27	1.09	1 (5%)	21,38,41	2.50	3 (14%)
1	UR3	0	2619	1	12,22,23	0.96	0	16,32,35	0.77	0
1	PSU	0	2621	1	13,21,22	1.56	2 (15%)	18,30,33	6.09	3 (16%)
1	1MA	0	628	1	14,25,26	1.00	1 (7%)	15,37,40	1.21	1 (6%)
3	5AA	4	76	1,3	16,26,27	0.65	0	15,38,41	1.22	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OMU	0	2587	1	-	0/5/27/28	0/2/2/2
1	OMG	0	2588	1,3	-	0/5/27/28	0/3/3/3
1	UR3	0	2619	1	-	0/3/25/26	0/2/2/2
1	PSU	0	2621	1	-	0/7/25/26	0/2/2/2
1	1MA	0	628	1	-	0/3/25/26	0/3/3/3
3	5AA	4	76	1,3	-	0/7/29/30	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	0	2621	PSU	C5-C1'	-4.49	1.48	1.52
1	0	2587	OMU	C4-N3	2.32	1.37	1.33
1	0	2621	PSU	C4-N3	2.75	1.38	1.33
1	0	628	1MA	C6-N6	2.78	1.34	1.29
1	0	2588	OMG	C6-N1	3.37	1.39	1.33

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0	2621	PSU	N1-C2-N3	-21.29	114.75	128.33
1	0	2588	OMG	C5-C6-N1	-8.61	111.81	123.59
1	0	628	1MA	C2-N3-C4	-3.62	110.80	116.40
1	0	2587	OMU	C5-C4-N3	-3.28	114.70	123.12
1	0	2588	OMG	N3-C2-N1	-2.32	123.91	127.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	0	2587	OMU	2	0
1	0	2619	UR3	1	0

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 233 ligands modelled in this entry, 233 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 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	0	2749/2922 (94%)	-0.49	34 (1%) 81 81	29, 51, 94, 161	0
2	9	122/122 (100%)	-0.46	4 (3%) 50 50	39, 62, 91, 152	0
3	4	5/8 (62%)	-0.44	0 100 100	41, 43, 47, 47	0
4	A	237/240 (98%)	0.18	10 (4%) 40 39	33, 57, 95, 119	0
5	B	337/338 (99%)	0.05	7 (2%) 67 68	29, 57, 80, 94	0
6	C	246/246 (100%)	-0.07	3 (1%) 81 81	27, 53, 75, 84	0
7	D	140/177 (79%)	2.21	63 (45%) 0 0	58, 101, 127, 137	0
8	E	172/178 (96%)	0.67	15 (8%) 13 10	47, 67, 87, 93	0
9	F	119/120 (99%)	0.96	22 (18%) 2 1	60, 82, 104, 119	0
10	G	29/348 (8%)	1.80	10 (34%) 0 0	65, 90, 102, 105	0
11	H	160/171 (93%)	0.30	8 (5%) 32 31	41, 59, 90, 99	0
12	J	142/145 (97%)	-0.01	1 (0%) 89 90	37, 52, 72, 90	0
13	K	132/132 (100%)	0.04	2 (1%) 76 76	34, 56, 77, 87	0
14	L	145/165 (87%)	0.68	24 (16%) 2 2	30, 72, 118, 132	0
15	M	194/194 (100%)	-0.08	1 (0%) 91 93	34, 50, 67, 74	0
16	N	186/187 (99%)	0.63	20 (10%) 8 6	38, 66, 115, 121	0
17	O	115/116 (99%)	0.13	0 100 100	45, 61, 77, 85	0
18	P	143/149 (95%)	0.29	0 100 100	45, 60, 77, 85	0
19	Q	95/96 (98%)	-0.03	1 (1%) 82 83	40, 48, 60, 75	0
20	R	150/155 (96%)	-0.15	1 (0%) 89 90	37, 48, 68, 75	0
21	S	81/85 (95%)	0.29	3 (3%) 45 45	49, 67, 85, 94	0
22	T	119/120 (99%)	0.75	13 (10%) 7 5	46, 63, 91, 110	0
23	U	53/66 (80%)	0.23	3 (5%) 27 26	43, 57, 74, 85	0
24	V	65/71 (91%)	1.79	23 (35%) 0 0	63, 86, 117, 121	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
25	W	154/154 (100%)	-0.04	2 (1%) 79 79	39, 50, 67, 76	0
26	X	82/92 (89%)	0.48	9 (10%) 7 5	45, 59, 77, 95	0
27	Y	142/241 (58%)	0.03	5 (3%) 48 48	30, 49, 71, 86	0
28	Z	73/83 (87%)	0.20	4 (5%) 29 27	44, 63, 77, 97	0
29	1	56/57 (98%)	-0.40	0 100 100	33, 40, 46, 54	0
30	2	46/50 (92%)	0.56	5 (10%) 7 5	41, 65, 97, 109	0
31	3	92/92 (100%)	0.33	3 (3%) 50 50	37, 59, 73, 84	0
32	I	70/162 (43%)	3.69	56 (80%) 0 0	99, 121, 143, 145	0
All	All	6651/7482 (88%)	0.01	352 (5%) 30 28	27, 56, 102, 161	0

The worst 5 of 352 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
24	V	1	THR	12.2
32	I	133	THR	10.4
32	I	71	GLY	10.2
7	D	63	ILE	9.5
32	I	93	GLN	8.7

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	OMG	0	2588	24/25	0.98	0.13	-	32,36,39,40	0
3	5AA	4	76	24/25	0.98	0.14	-	38,44,48,48	0
1	1MA	0	628	23/24	0.99	0.15	-	29,32,34,36	0
1	UR3	0	2619	21/22	0.99	0.15	-	32,39,41,42	0
1	PSU	0	2621	20/21	0.99	0.14	-	30,33,37,38	0
1	OMU	0	2587	21/22	0.99	0.13	-	35,37,39,42	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
36	CL	0	9315	1/1	0.93	0.41	65.34	84,84,84,84	0
35	NA	0	9174	1/1	0.86	0.89	41.97	64,64,64,64	0
35	NA	0	9171	1/1	0.73	0.41	31.45	64,64,64,64	0
35	NA	R	9186	1/1	0.30	0.75	22.47	84,84,84,84	0
35	NA	0	9156	1/1	0.95	0.45	21.62	51,51,51,51	0
35	NA	0	9177	1/1	0.80	0.41	18.69	73,73,73,73	0
35	NA	0	9182	1/1	0.73	0.45	18.48	84,84,84,84	0
35	NA	L	9180	1/1	0.93	0.52	17.97	72,72,72,72	0
35	NA	0	9150	1/1	0.88	0.29	15.99	48,48,48,48	0
35	NA	0	9135	1/1	0.93	0.28	15.22	50,50,50,50	0
35	NA	0	9125	1/1	0.93	0.26	15.15	69,69,69,69	0
35	NA	0	9172	1/1	0.93	0.38	12.80	61,61,61,61	0
35	NA	0	9155	1/1	0.92	0.54	12.49	78,78,78,78	0
33	MG	0	8060	1/1	0.99	0.31	11.35	49,49,49,49	0
35	NA	0	9161	1/1	0.90	0.30	7.75	56,56,56,56	0
35	NA	0	9121	1/1	0.96	0.33	7.27	58,58,58,58	0
35	NA	0	9178	1/1	0.97	0.21	6.37	57,57,57,57	0
35	NA	0	9162	1/1	0.95	0.23	6.02	72,72,72,72	0
34	K	0	9001	1/1	0.96	0.25	5.43	74,74,74,74	0
35	NA	0	9129	1/1	0.69	0.17	4.72	65,65,65,65	0
35	NA	0	9168	1/1	0.84	0.15	3.77	56,56,56,56	0
35	NA	0	9173	1/1	0.90	0.24	3.45	58,58,58,58	0
35	NA	0	9102	1/1	0.96	0.16	3.26	44,44,44,44	0
33	MG	0	8053	1/1	0.98	0.17	3.13	58,58,58,58	0
35	NA	Q	9148	1/1	0.93	0.29	2.96	43,43,43,43	0
35	NA	M	9147	1/1	0.98	0.21	2.31	33,33,33,33	0
35	NA	0	9126	1/1	0.85	0.18	2.16	44,44,44,44	0
35	NA	0	9176	1/1	0.97	0.18	1.84	42,42,42,42	0
33	MG	0	8080	1/1	0.95	0.15	1.73	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
36	CL	0	9305	1/1	0.85	0.18	1.60	71,71,71,71	0
33	MG	0	8054	1/1	0.91	0.15	1.47	39,39,39,39	0
33	MG	0	8013	1/1	0.98	0.19	1.45	41,41,41,41	0
35	NA	0	9110	1/1	0.92	0.16	1.29	38,38,38,38	0
33	MG	A	8065	1/1	0.98	0.21	1.22	52,52,52,52	0
33	MG	0	8038	1/1	0.99	0.16	1.06	26,26,26,26	0
33	MG	0	8012	1/1	0.97	0.16	1.04	32,32,32,32	0
35	NA	0	9124	1/1	0.86	0.17	0.83	67,67,67,67	0
36	CL	0	9316	1/1	0.96	0.16	0.62	61,61,61,61	0
33	MG	0	8018	1/1	0.99	0.15	0.47	47,47,47,47	0
36	CL	O	9308	1/1	0.97	0.22	0.40	81,81,81,81	0
35	NA	0	9103	1/1	0.97	0.15	0.35	39,39,39,39	0
35	NA	0	9165	1/1	0.93	0.21	0.22	42,42,42,42	0
35	NA	9	9183	1/1	0.89	0.13	-0.23	49,49,49,49	0
35	NA	A	9145	1/1	0.96	0.17	-0.28	45,45,45,45	0
36	CL	0	9313	1/1	0.98	0.14	-0.28	61,61,61,61	0
33	MG	0	8044	1/1	0.94	0.12	-0.38	49,49,49,49	0
35	NA	R	9137	1/1	0.85	0.12	-0.46	50,50,50,50	0
36	CL	L	9310	1/1	0.98	0.15	-0.48	61,61,61,61	0
35	NA	H	9109	1/1	0.95	0.13	-0.60	36,36,36,36	0
35	NA	C	9104	1/1	0.89	0.15	-0.62	41,41,41,41	0
33	MG	B	8055	1/1	0.98	0.14	-0.67	52,52,52,52	0
36	CL	M	9318	1/1	0.98	0.15	-0.78	50,50,50,50	0
33	MG	0	8077	1/1	0.99	0.13	-0.82	35,35,35,35	0
35	NA	0	9123	1/1	0.96	0.17	-0.86	43,43,43,43	0
33	MG	0	8057	1/1	0.95	0.14	-0.93	45,45,45,45	0
35	NA	0	9166	1/1	0.84	0.11	-0.96	68,68,68,68	0
35	NA	0	9133	1/1	0.92	0.11	-1.00	32,32,32,32	0
33	MG	0	8107	1/1	0.98	0.09	-1.06	38,38,38,38	0
35	NA	0	9144	1/1	0.97	0.10	-1.08	27,27,27,27	0
33	MG	0	8033	1/1	0.99	0.13	-1.11	32,32,32,32	0
36	CL	J	9321	1/1	0.97	0.13	-1.12	54,54,54,54	0
35	NA	0	9114	1/1	0.96	0.12	-1.18	51,51,51,51	0
37	CD	Z	9203	1/1	0.99	0.09	-1.21	68,68,68,68	0
37	CD	U	9201	1/1	0.99	0.09	-1.21	69,69,69,69	0
33	MG	0	8017	1/1	0.98	0.10	-1.25	33,33,33,33	0
33	MG	0	8064	1/1	0.96	0.12	-1.27	33,33,33,33	0
36	CL	B	9319	1/1	0.99	0.14	-1.37	60,60,60,60	0
35	NA	0	9117	1/1	0.92	0.11	-1.44	46,46,46,46	0
33	MG	0	8015	1/1	0.99	0.11	-1.48	31,31,31,31	0
33	MG	0	8086	1/1	0.97	0.06	-1.61	47,47,47,47	0
35	NA	0	9139	1/1	0.99	0.13	-1.75	22,22,22,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
33	MG	0	8058	1/1	0.98	0.12	-1.76	57,57,57,57	0
33	MG	0	8074	1/1	0.99	0.09	-1.82	40,40,40,40	0
33	MG	0	8101	1/1	0.92	0.11	-1.86	74,74,74,74	0
36	CL	3	9304	1/1	0.93	0.15	-1.96	70,70,70,70	0
33	MG	T	8073	1/1	0.90	0.09	-1.97	66,66,66,66	0
36	CL	K	9312	1/1	0.97	0.10	-1.98	55,55,55,55	0
37	CD	1	9202	1/1	1.00	0.05	-2.07	68,68,68,68	0
33	MG	0	8112	1/1	0.95	0.11	-2.17	47,47,47,47	0
35	NA	J	9146	1/1	0.98	0.10	-2.22	35,35,35,35	0
33	MG	0	8056	1/1	0.97	0.12	-2.42	51,51,51,51	0
35	NA	0	9127	1/1	0.95	0.10	-2.44	35,35,35,35	0
33	MG	0	8070	1/1	0.98	0.04	-2.47	45,45,45,45	0
33	MG	0	8062	1/1	0.96	0.10	-2.66	57,57,57,57	0
33	MG	0	8003	1/1	0.99	0.14	-2.68	40,40,40,40	0
33	MG	0	8108	1/1	0.97	0.10	-2.71	75,75,75,75	0
33	MG	0	8008	1/1	0.99	0.08	-2.82	37,37,37,37	0
33	MG	3	8078	1/1	0.93	0.05	-2.89	55,55,55,55	0
33	MG	0	8010	1/1	0.99	0.11	-2.99	29,29,29,29	0
35	NA	0	9131	1/1	0.97	0.08	-3.01	36,36,36,36	0
33	MG	0	8020	1/1	0.99	0.12	-3.27	24,24,24,24	0
33	MG	0	8096	1/1	0.96	0.07	-3.33	51,51,51,51	0
35	NA	R	9138	1/1	0.91	0.06	-3.47	61,61,61,61	0
35	NA	0	9153	1/1	0.98	0.10	-3.56	25,25,25,25	0
35	NA	0	9132	1/1	0.94	0.09	-3.71	35,35,35,35	0
33	MG	0	8001	1/1	0.99	0.10	-3.74	38,38,38,38	0
33	MG	0	8019	1/1	0.99	0.09	-3.95	38,38,38,38	0
33	MG	0	8004	1/1	0.98	0.05	-3.97	27,27,27,27	0
33	MG	0	8091	1/1	0.99	0.07	-4.38	57,57,57,57	0
33	MG	0	8007	1/1	0.99	0.07	-4.69	24,24,24,24	0
33	MG	0	8032	1/1	0.98	0.05	-4.78	35,35,35,35	0
33	MG	0	8022	1/1	0.98	0.07	-4.85	35,35,35,35	0
33	MG	0	8084	1/1	0.94	0.10	-5.05	51,51,51,51	0
33	MG	0	8067	1/1	0.98	0.09	-5.20	46,46,46,46	0
33	MG	0	8118	1/1	0.99	0.09	-5.30	40,40,40,40	0
34	K	0	9002	1/1	0.98	0.08	-5.34	54,54,54,54	0
33	MG	0	8006	1/1	0.99	0.05	-5.51	33,33,33,33	0
33	MG	Y	8109	1/1	0.99	0.09	-5.73	39,39,39,39	0
35	NA	0	9120	1/1	0.99	0.10	-5.81	43,43,43,43	0
35	NA	0	9105	1/1	0.97	0.08	-6.61	39,39,39,39	0
33	MG	0	8035	1/1	0.98	0.07	-6.80	51,51,51,51	0
35	NA	0	9143	1/1	0.98	0.07	-8.25	36,36,36,36	0
33	MG	0	8002	1/1	0.99	0.05	-8.54	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
33	MG	0	8110	1/1	0.99	0.07	-10.22	40,40,40,40	0
33	MG	9	8052	1/1	0.98	0.05	-12.56	47,47,47,47	0
33	MG	0	8023	1/1	0.98	0.18	-	48,48,48,48	0
33	MG	0	8029	1/1	0.99	0.17	-	36,36,36,36	0
33	MG	0	8117	1/1	0.98	0.10	-	40,40,40,40	0
35	NA	0	9185	1/1	0.90	0.65	-	54,54,54,54	0
33	MG	0	8102	1/1	0.97	0.14	-	67,67,67,67	0
33	MG	0	8104	1/1	0.96	0.21	-	73,73,73,73	0
35	NA	0	9163	1/1	0.91	0.41	-	63,63,63,63	0
35	NA	0	9128	1/1	0.97	0.07	-	35,35,35,35	0
33	MG	0	8094	1/1	0.95	0.12	-	86,86,86,86	0
33	MG	0	8089	1/1	0.92	0.09	-	60,60,60,60	0
33	MG	0	8011	1/1	0.98	0.16	-	23,23,23,23	0
35	NA	0	9152	1/1	0.86	0.41	-	65,65,65,65	0
35	NA	0	9136	1/1	0.98	0.07	-	56,56,56,56	0
35	NA	0	9184	1/1	0.94	0.53	-	96,96,96,96	0
36	CL	0	9322	1/1	0.83	0.54	-	92,92,92,92	0
33	MG	9	8095	1/1	0.85	0.14	-	75,75,75,75	0
33	MG	0	8100	1/1	0.99	0.18	-	82,82,82,82	0
37	CD	3	9204	1/1	0.95	0.09	-	66,66,66,66	0
35	NA	0	9149	1/1	0.90	0.14	-	42,42,42,42	0
33	MG	0	8043	1/1	0.91	0.08	-	53,53,53,53	0
33	MG	0	8082	1/1	0.86	0.20	-	65,65,65,65	0
33	MG	0	8114	1/1	0.94	0.18	-	56,56,56,56	0
35	NA	0	9160	1/1	0.94	0.33	-	49,49,49,49	0
35	NA	0	9181	1/1	0.95	0.12	-	48,48,48,48	0
33	MG	0	8016	1/1	0.98	0.21	-	44,44,44,44	0
33	MG	0	8045	1/1	0.97	0.11	-	62,62,62,62	0
33	MG	0	8115	1/1	0.96	0.12	-	49,49,49,49	0
34	K	0	9003	1/1	0.90	0.13	-	66,66,66,66	0
33	MG	0	8093	1/1	0.84	0.11	-	52,52,52,52	0
33	MG	0	8081	1/1	0.96	0.18	-	54,54,54,54	0
35	NA	0	9101	1/1	0.92	0.21	-	47,47,47,47	0
35	NA	0	9167	1/1	0.93	0.08	-	47,47,47,47	0
33	MG	0	8090	1/1	0.76	0.60	-	69,69,69,69	0
33	MG	0	8088	1/1	0.93	0.06	-	36,36,36,36	0
36	CL	0	9317	1/1	0.97	0.08	-	64,64,64,64	0
36	CL	0	9311	1/1	0.99	0.09	-	52,52,52,52	0
36	CL	J	9302	1/1	0.94	0.06	-	61,61,61,61	0
35	NA	0	9164	1/1	0.92	0.23	-	55,55,55,55	0
35	NA	0	9106	1/1	0.95	0.80	-	51,51,51,51	0
33	MG	0	8026	1/1	0.99	0.13	-	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
33	MG	0	8009	1/1	0.99	0.11	-	35,35,35,35	0
35	NA	0	9157	1/1	0.95	0.09	-	69,69,69,69	0
33	MG	0	8014	1/1	0.95	0.11	-	41,41,41,41	0
33	MG	0	8042	1/1	0.94	0.08	-	39,39,39,39	0
35	NA	0	9113	1/1	0.94	0.16	-	66,66,66,66	0
36	CL	R	9306	1/1	0.96	0.12	-	48,48,48,48	0
33	MG	0	8025	1/1	0.99	0.11	-	35,35,35,35	0
33	MG	0	8027	1/1	0.95	0.13	-	55,55,55,55	0
33	MG	0	8036	1/1	0.99	0.13	-	35,35,35,35	0
33	MG	0	8103	1/1	0.97	0.11	-	81,81,81,81	0
35	NA	0	9130	1/1	0.98	0.08	-	44,44,44,44	0
36	CL	J	9301	1/1	0.98	0.19	-	73,73,73,73	0
33	MG	0	8113	1/1	0.96	0.12	-	45,45,45,45	0
33	MG	0	8051	1/1	0.96	0.07	-	61,61,61,61	0
33	MG	0	8034	1/1	0.98	0.10	-	32,32,32,32	0
33	MG	0	8068	1/1	0.96	0.08	-	59,59,59,59	0
33	MG	0	8021	1/1	0.99	0.17	-	30,30,30,30	0
35	NA	0	9154	1/1	0.97	0.19	-	38,38,38,38	0
33	MG	0	8111	1/1	0.94	0.15	-	77,77,77,77	0
33	MG	0	8031	1/1	1.00	0.08	-	37,37,37,37	0
35	NA	0	9175	1/1	0.97	0.45	-	56,56,56,56	0
33	MG	0	8071	1/1	0.94	0.13	-	72,72,72,72	0
36	CL	0	9320	1/1	0.96	0.11	-	48,48,48,48	0
36	CL	N	9307	1/1	0.94	0.20	-	66,66,66,66	0
35	NA	0	9111	1/1	0.92	0.11	-	59,59,59,59	0
35	NA	0	9169	1/1	0.93	0.38	-	79,79,79,79	0
33	MG	K	8069	1/1	0.84	0.15	-	49,49,49,49	0
33	MG	0	8075	1/1	0.96	0.06	-	43,43,43,43	0
35	NA	0	9159	1/1	0.94	0.23	-	56,56,56,56	0
35	NA	0	9108	1/1	0.97	0.09	-	61,61,61,61	0
33	MG	0	8116	1/1	0.97	0.06	-	47,47,47,47	0
33	MG	0	8085	1/1	0.80	0.24	-	74,74,74,74	0
36	CL	A	9309	1/1	0.89	0.19	-	77,77,77,77	0
33	MG	0	8040	1/1	0.97	0.18	-	52,52,52,52	0
35	NA	0	9158	1/1	0.90	0.59	-	84,84,84,84	0
33	MG	0	8049	1/1	0.87	0.29	-	80,80,80,80	0
33	MG	0	8106	1/1	0.91	0.08	-	63,63,63,63	0
36	CL	0	9303	1/1	0.94	0.13	-	64,64,64,64	0
33	MG	0	8076	1/1	0.93	0.06	-	70,70,70,70	0
36	CL	0	9314	1/1	0.97	0.07	-	53,53,53,53	0
33	MG	0	8092	1/1	0.72	0.66	-	115,115,115,115	0
35	NA	S	9112	1/1	0.86	0.51	-	75,75,75,75	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
35	NA	0	9141	1/1	0.96	0.07	-	47,47,47,47	0
35	NA	0	9142	1/1	0.91	0.17	-	53,53,53,53	0
33	MG	0	8046	1/1	0.96	0.09	-	54,54,54,54	0
33	MG	0	8028	1/1	0.97	0.12	-	44,44,44,44	0
33	MG	0	8059	1/1	0.96	0.07	-	44,44,44,44	0
33	MG	0	8005	1/1	1.00	0.14	-	37,37,37,37	0
33	MG	A	8066	1/1	0.92	0.05	-	66,66,66,66	0
35	NA	0	9140	1/1	0.94	0.20	-	45,45,45,45	0
35	NA	0	9134	1/1	0.98	0.10	-	37,37,37,37	0
33	MG	0	8048	1/1	0.97	0.16	-	62,62,62,62	0
33	MG	0	8087	1/1	0.95	0.14	-	54,54,54,54	0
33	MG	0	8047	1/1	0.92	0.18	-	78,78,78,78	0
33	MG	0	8039	1/1	0.93	0.06	-	51,51,51,51	0
35	NA	0	9115	1/1	0.96	0.15	-	39,39,39,39	0
33	MG	0	8072	1/1	0.99	0.24	-	55,55,55,55	0
37	CD	O	9205	1/1	0.04	0.30	-	200,200,200,200	0
33	MG	0	8030	1/1	1.00	0.15	-	26,26,26,26	0
33	MG	0	8061	1/1	0.98	0.08	-	45,45,45,45	0
35	NA	0	9118	1/1	0.98	0.20	-	47,47,47,47	0
33	MG	0	8041	1/1	0.90	0.24	-	72,72,72,72	0
35	NA	0	9116	1/1	0.95	0.11	-	36,36,36,36	0
33	MG	0	8079	1/1	0.99	0.12	-	38,38,38,38	0
35	NA	9	9151	1/1	0.63	0.36	-	87,87,87,87	0
33	MG	0	8099	1/1	0.95	0.10	-	63,63,63,63	0
33	MG	0	8050	1/1	0.91	0.08	-	61,61,61,61	0
35	NA	0	9170	1/1	0.85	0.48	-	97,97,97,97	0
33	MG	0	8063	1/1	0.95	0.15	-	84,84,84,84	0
33	MG	0	8097	1/1	0.95	0.15	-	40,40,40,40	0
33	MG	0	8083	1/1	0.94	0.08	-	42,42,42,42	0
33	MG	0	8024	1/1	0.83	0.65	-	83,83,83,83	0
35	NA	0	9119	1/1	0.98	0.08	-	36,36,36,36	0
33	MG	0	8037	1/1	0.98	0.06	-	44,44,44,44	0
35	NA	0	9179	1/1	0.95	0.18	-	57,57,57,57	0
35	NA	0	9107	1/1	0.91	0.20	-	50,50,50,50	0
33	MG	0	8098	1/1	0.93	0.24	-	40,40,40,40	0
35	NA	H	9122	1/1	0.96	0.13	-	59,59,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.