



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 11:17 am BST

PDB ID : 1NJI  
Title : Structure of chloramphenicol bound to the 50S ribosomal subunit  
Authors : Hansen, J.L.; Moore, P.B.; Steitz, T.A.  
Deposited on : 2002-12-31  
Resolution : 3.00 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

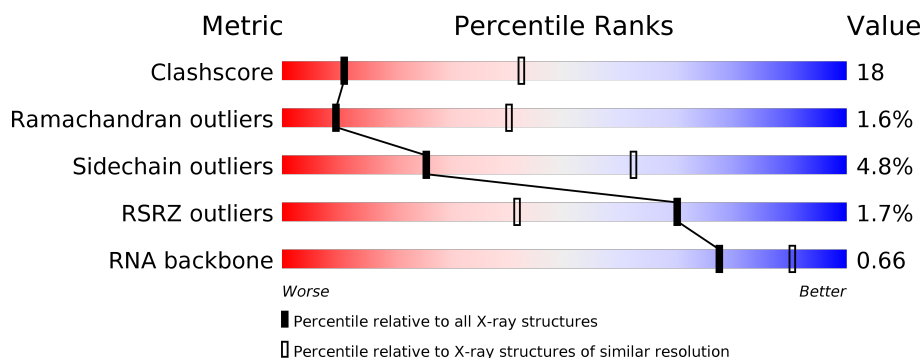
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



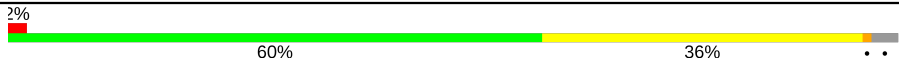

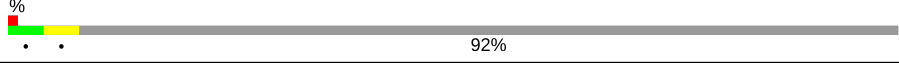
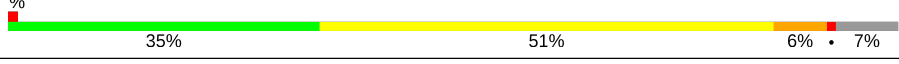

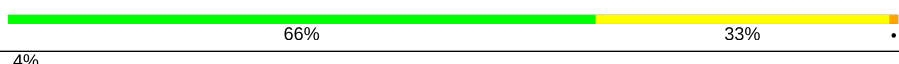
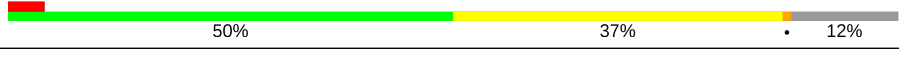
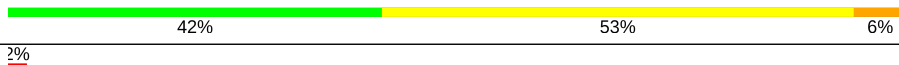
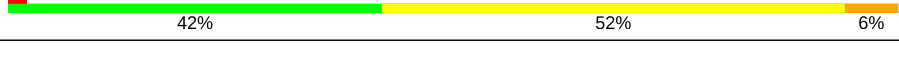


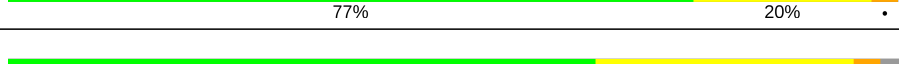





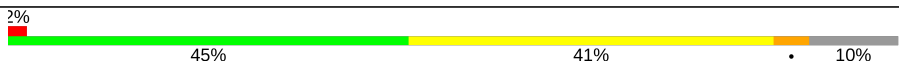
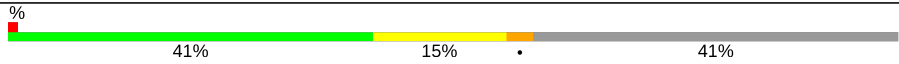



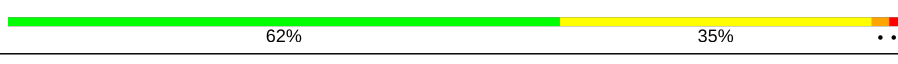

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)
RNA backbone	3102	1173 (3.30-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria 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	<div> <div>57%</div> <div>31%</div> <div>6%</div> <div>6%</div> </div>
2	B	122	<div>5%</div> <div>56%</div> <div>30%</div> <div>11%</div> <div>•</div>
3	C	239	<div>54%</div> <div>39%</div> <div>6%</div> <div>•</div>
4	D	337	<div>51%</div> <div>44%</div> <div>5%</div>
5	E	246	<div>57%</div> <div>38%</div> <div>5%</div>
6	F	176	<div>18%</div> <div>28%</div> <div>43%</div> <div>7%</div> <div>•</div> <div>20%</div>

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Mol	Chain	Length	Quality of chain
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	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
31	MG	A	8049	-	-	-	X
33	NA	A	8371	-	-	-	X
33	NA	A	8376	-	-	-	X
33	NA	A	8377	-	-	-	X
33	NA	A	8382	-	-	-	X
33	NA	A	8384	-	-	-	X

## 2 Entry composition

There are 37 unique types of molecules in this entry. The entry contains 98536 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	A	2754	Total	C	N	O	P	0	0	0
			59017	26346	10878	19048	2745			

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

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 50S ribosomal protein L2P.

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 50S ribosomal protein L3P.

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	SEE REMARK 999	UNP P20279
D	310	ARG	PHE	SEE REMARK 999	UNP P20279

- Molecule 5 is a protein called 50S ribosomal protein L4E.

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 50S ribosomal protein L5P.

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 50S ribosomal protein L6P.

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 50S 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 Acidic ribosomal protein P0 homolog.

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 50S 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 50S ribosomal protein L13P.

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 50S ribosomal protein L14P.

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 50S ribosomal protein L15P.

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

- Molecule 14 is a protein called 50S 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 50S ribosomal protein L18P.

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 50S ribosomal protein L18E.

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

- Molecule 17 is a protein called 50S ribosomal protein L19E.

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

- Molecule 18 is a protein called 50S ribosomal protein L21e.

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

- Molecule 19 is a protein called 50S ribosomal protein L22P.

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

- Molecule 20 is a protein called 50S ribosomal protein L23P.

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

- Molecule 21 is a protein called 50S ribosomal protein L24P.

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

- Molecule 22 is a protein called 50S ribosomal protein L24E.

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

- Molecule 23 is a protein called 50S ribosomal protein L29P.

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

- Molecule 24 is a protein called 50S ribosomal protein L30P.

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

- Molecule 25 is a protein called 50S 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 50S 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 50S 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 50S 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 50S 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			

- Molecule 30 is a protein called 50S 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 MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
31	1	1	Total	Mg	0	0
			1	1		
31	D	1	Total	Mg	0	0
			1	1		
31	B	1	Total	Mg	0	0
			1	1		
31	C	1	Total	Mg	0	0
			1	1		
31	Z	1	Total	Mg	0	0
			1	1		
31	A	109	Total	Mg	0	0
			109	109		
31	4	1	Total	Mg	0	0
			1	1		
31	U	1	Total	Mg	0	0
			1	1		
31	L	1	Total	Mg	0	0
			1	1		

- Molecule 32 is POTASSIUM ION (three-letter code: K) (formula: K).

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

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

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

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

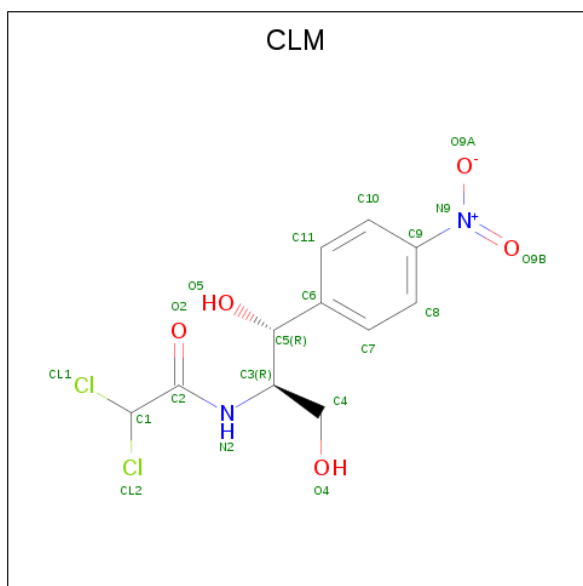
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
34	P	1	Total Cl 1 1	0	0
34	D	1	Total Cl 1 1	0	0
34	K	3	Total Cl 3 3	0	0
34	C	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
34	Z	1	Total	Cl	0	0
			1	1		
34	A	9	Total	Cl	0	0
			9	9		
34	4	1	Total	Cl	0	0
			1	1		
34	N	1	Total	Cl	0	0
			1	1		
34	O	1	Total	Cl	0	0
			1	1		
34	R	1	Total	Cl	0	0
			1	1		
34	S	1	Total	Cl	0	0
			1	1		
34	M	1	Total	Cl	0	0
			1	1		

- Molecule 35 is CHLORAMPHENICOL (three-letter code: CLM) (formula:  $C_{11}H_{12}Cl_2N_2O_5$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
35	A	1	Total	C	Cl	N	O	
			20	11	2	2	5	

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

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

- Molecule 37 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
37	A	5906	Total O 5906 5906	0	0
37	B	143	Total O 143 143	0	0
37	C	123	Total O 123 123	0	0
37	D	147	Total O 147 147	0	0
37	E	167	Total O 167 167	0	0
37	F	50	Total O 50 50	0	0
37	G	45	Total O 45 45	0	0
37	H	29	Total O 29 29	0	0
37	I	21	Total O 21 21	0	0
37	J	76	Total O 76 76	0	0
37	K	54	Total O 54 54	0	0
37	L	61	Total O 61 61	0	0
37	M	78	Total O 78 78	0	0
37	N	125	Total O 125 125	0	0
37	O	69	Total O 69 69	0	0

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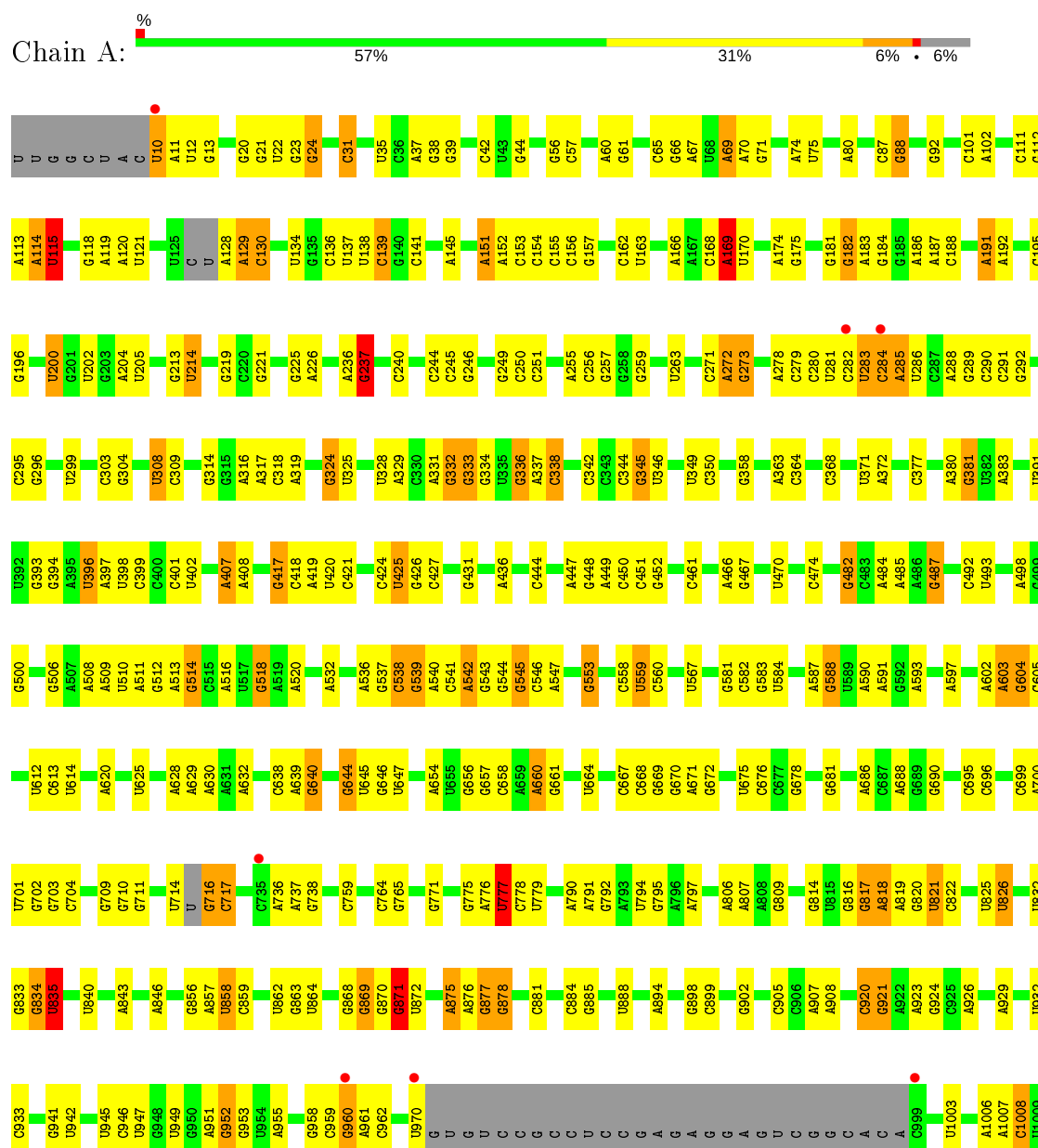
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	P	45	Total 45	O 45	0	0
37	Q	73	Total 73	O 73	0	0
37	R	53	Total 53	O 53	0	0
37	S	85	Total 85	O 85	0	0
37	T	35	Total 35	O 35	0	0
37	U	40	Total 40	O 40	0	0
37	V	28	Total 28	O 28	0	0
37	W	17	Total 17	O 17	0	0
37	X	69	Total 69	O 69	0	0
37	Y	28	Total 28	O 28	0	0
37	Z	98	Total 98	O 98	0	0
37	1	37	Total 37	O 37	0	0
37	2	54	Total 54	O 54	0	0
37	3	43	Total 43	O 43	0	0
37	4	68	Total 68	O 68	0	0

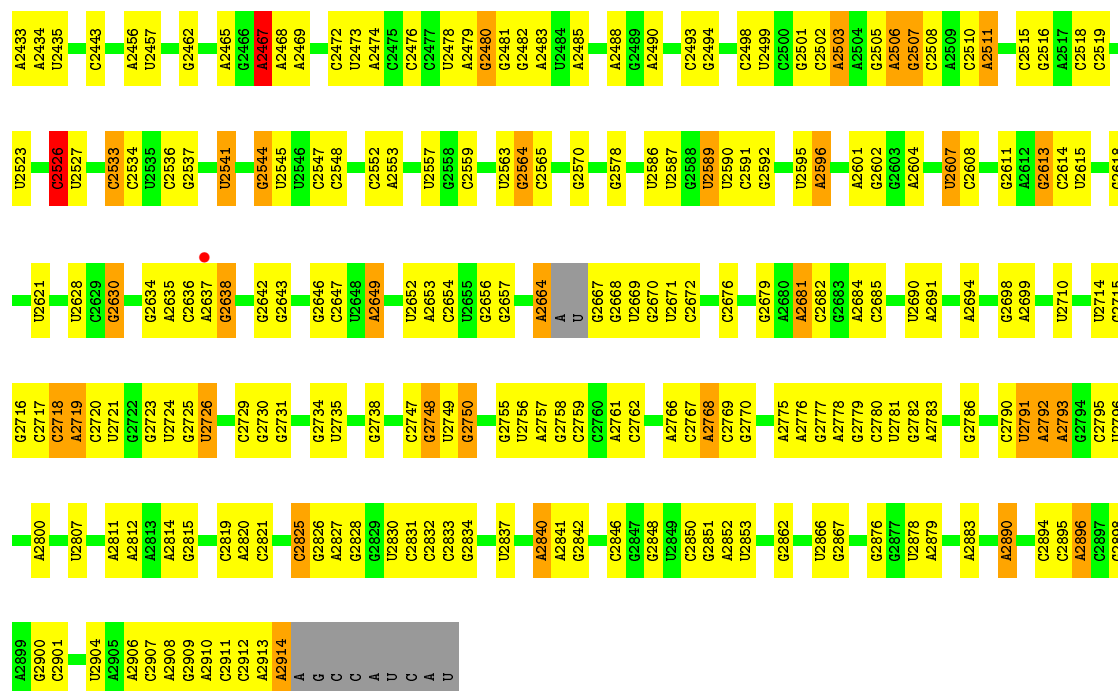
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

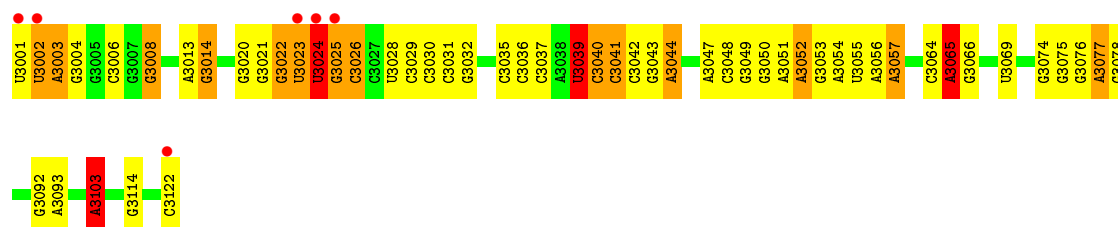
#### • Molecule 1: 23S ribosomal RNA



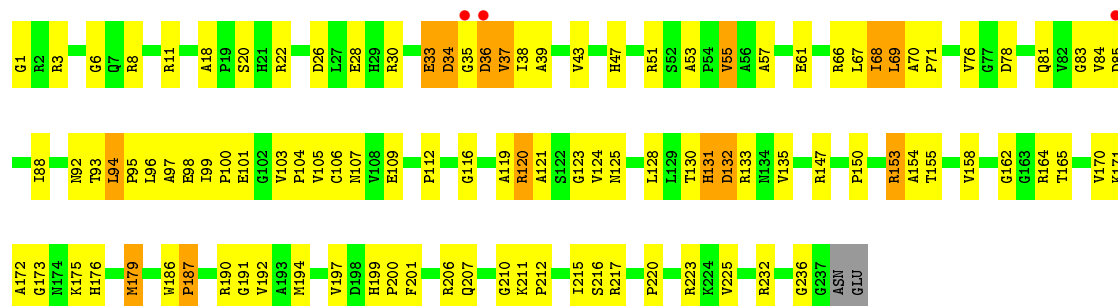
A2338	G2102	U1874	A1856	C1536	A1434	A1328	A1207	A1132	G1010
A2341	A2103	G1877	A1657	G1543	U1435	A1329	C1208	A1133	A1013
U2242	C2104	G1878	A1658	U1544	U1436	U1333	G1209	G1134	A1014
C2247	G2110	U1879	C1666	G1545	A1437	C1334	G1210	G1135	C1016
G2251	C2114	U1883	U1667	G1546	A1438	C1335	G1211	U1136	U1016
A2252	C2119	G1884	U1668	G1555	U1439	G1340	C1212	U1137	U1017
G2256	U2120	U1887	U1669	U1559	U1440	A1341	G1213	U1138	
G2257	G2121	U1887	G1670	A1559	U1441	C1342	A1215	U1139	U1028
A2258	C2122	U1887	C1671	U	A1442	C1343	G1216	U1140	U1029
	G2128	G1902	C1679	U	G1445		G1217	G1151	
U2265	A2010	G1903	G1680	U1561	U1446	G1351	U1218	G1044	G1045
A2266	A2011	A1904	G1681	C1562	U1447	A1352	U1219		
G2267	U2133	A1909	A1682	G1563		G1353	C1229	C1051	
G2268	G2134	A1910	G1683	C1564	C1450	C1354	U1234	G1052	
C2269	A2135	U1887	A1685	A1573	C1451	A1355	G1235	G1053	
G2270	G2136	U1788	G1688	A1580	U1457	C1360	U1236	U1056	
G2271	A	G1789	C1692	C1584	A1458	C1361	U1237	A1057	
G2272	C	C1798	C1692	C1585		U1362	U1238	A1058	
G2273	G	A1804	C1700	C1585	C1462	A1372	G1239	G1059	
A2274	U	G1805	A1701	G1588	A1463		A1242	C1060	
G2275	G	G1806	U1702	G1589	A1470	A1375	C1243	A1067	
U2276	C	U1933	G1706	G1592	C1474	G1376	U1244		
U2277	C	A1815	G1707	C1592		C1377	A1246	G1072	
G2281	G	C1818	C1593	C1594	U1477	U1383	C1245	G1076	
U2282	C	G1819	C1594	C1595	U1478	C1384	U1249	G1077	
G2283	A	A1942	A1717	U1596	C1483	G1391	C1250	A1078	
	C	C1943	U1722	A1597	G1484	A1392	C1251	A1079	
C2290	U	U1946	G1723	U1599	A1485	A1393	U1266	C1080	
A2291	G	G1951	U1725	A1603	A1494	C1394	C1267	A1081	
U2297	G	U	G1730	G1604	C1495	G1398	U1268	C1084	
A2300	C	A	C1731	G1605	G1496	A1399	C1269	C1085	
A2301	A	A	A1732	G1605	G1497		U1270	C1086	
A2302	C	U	C1733	C1613	U1500	A1406	C1273	G1087	
U2308	A	U	C1734	A1614	U1500	U1407	U1279	A1088	
G2310	C	U	C1735	A1615	U1503	G1409		A1098	
	C	A	A1736	U1625	A1504		C1289	G1099	
C2313	C	C	A1737	A1626	U1505	A1413	G1290	U1109	
G2314	C	C	C1738	G1627	U1506	A1414	U1298	G1110	
U2315	C	U	U1740			G1415	U1299		
G2316	U	U1964	U1741	A1630	A1515	G1417	G1299	A1114	
C2317	C	G1970	A1742	A1633	U1517	U1418	U1304	U1115	
	G	G1971	G1743	G1634	G1523	U1419	C1305	U1116	
U2320	A	U1972	G1751	U1635	U1524	C1420	U1306	A1117	
A2321	C	A1973	G1752	G1636	U1525	C1421	U1311	A1118	
	C	G1974	C1753	A1637	G1526	C1422	G1312	G1119	
G2324	A	U1978	A1754	A1641	A1527	C1423	A1313	U1120	
C2325	C	A1978	A1755	A1642	A1528	A1424	U1314	A1123	
U2326	U	G1979	G1756	A1642	G1529	G1425		C1201	
A2329	C	U1980	U1757	U1654		C1426	G1325	G1203	
U2330	C	A1981	U1758	G1655		G1430		U1205	
		C1982	A1759					U1300	
								G1206	



• Molecule 2: 5S ribosomal RNA



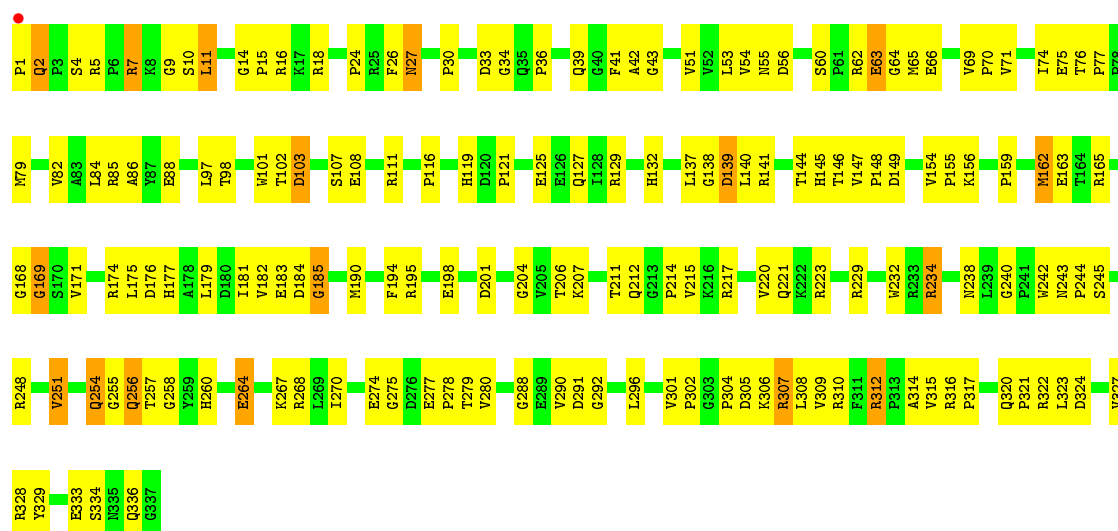
• Molecule 3: 50S ribosomal protein L2P



• Molecule 4: 50S ribosomal protein L3P

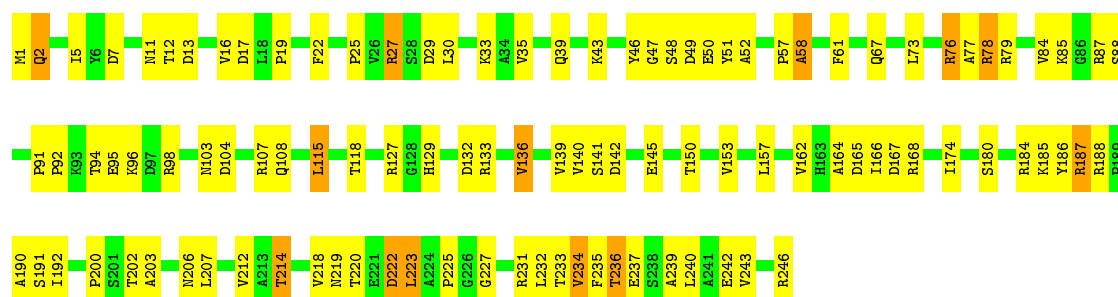






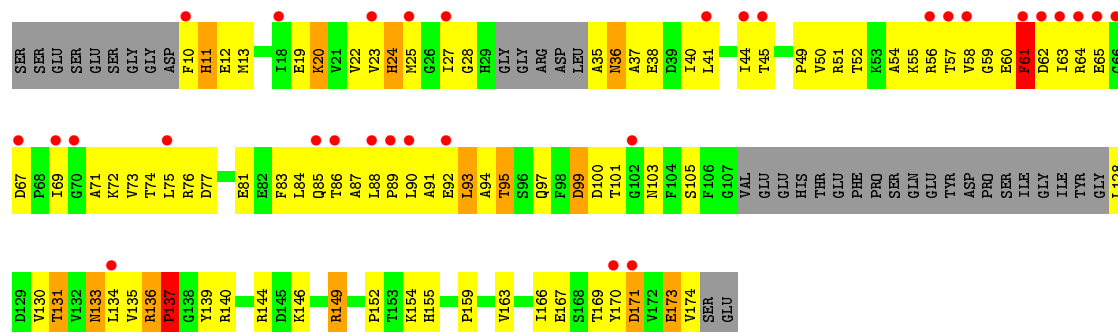
• Molecule 5: 50S ribosomal protein L4E

Chain E: 57% 38% 5%



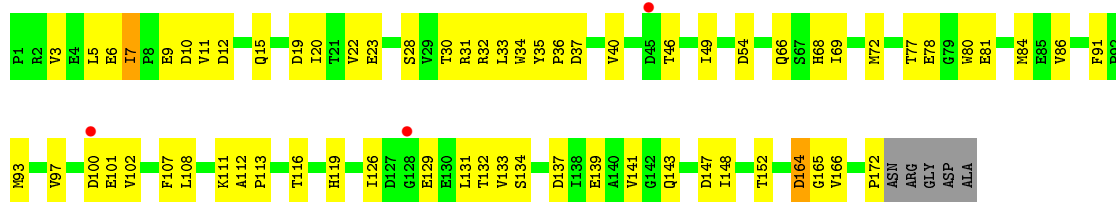
• Molecule 6: 50S ribosomal protein L5P

Chain F: 18% 28% 43% 7% 20%

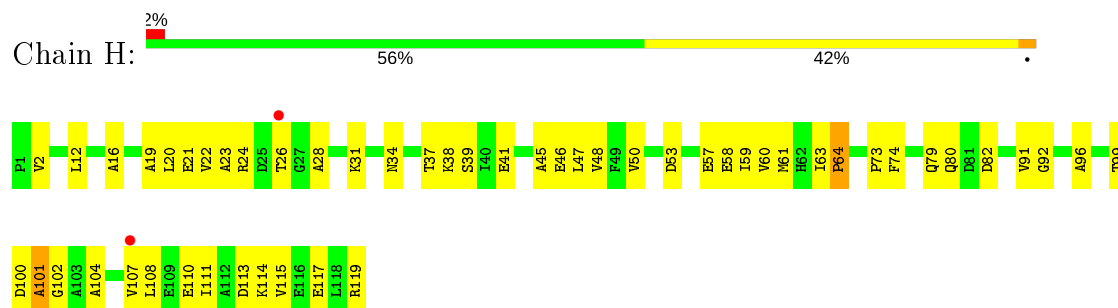


• Molecule 7: 50S ribosomal protein L6P

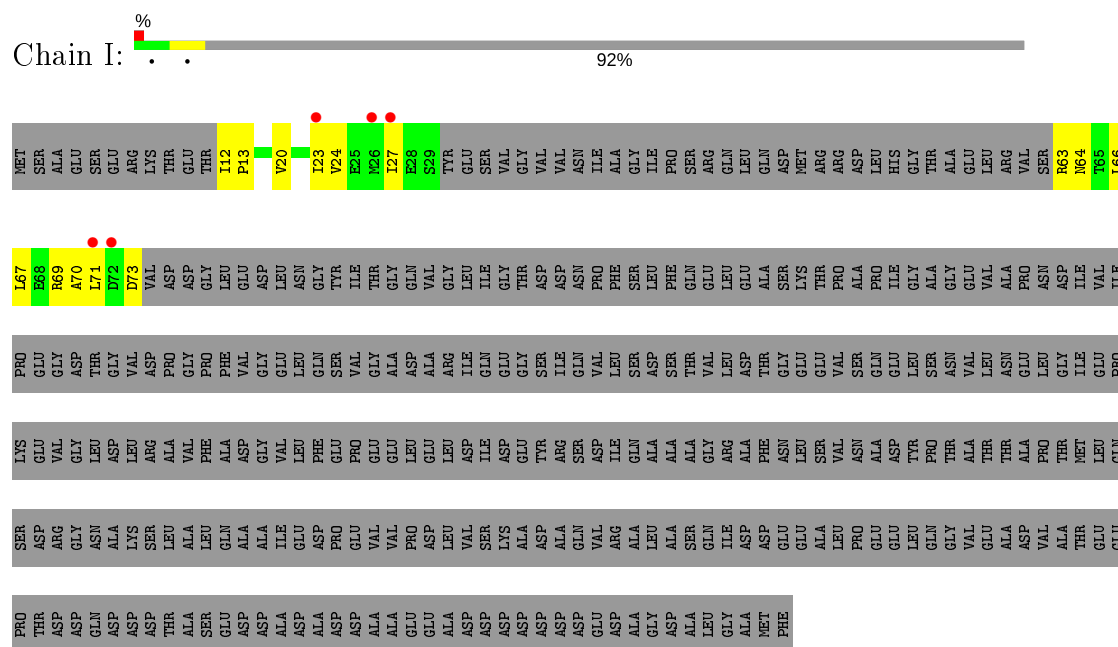
Chain G: 2% 60% 36%



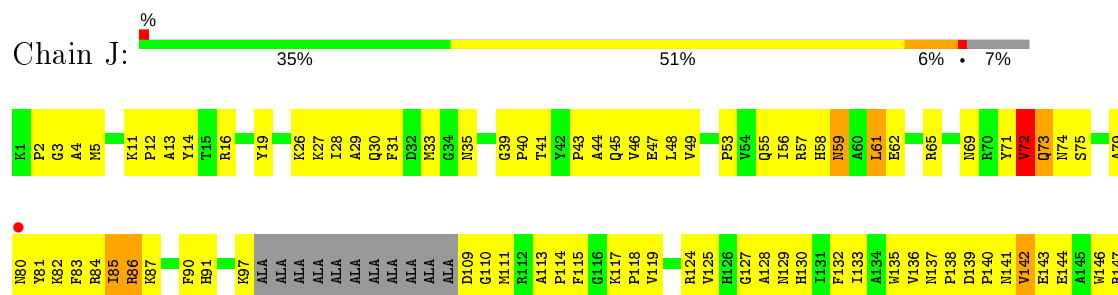
• Molecule 8: 50S ribosomal protein L7Ae



• Molecule 9: Acidic ribosomal protein P0 homolog



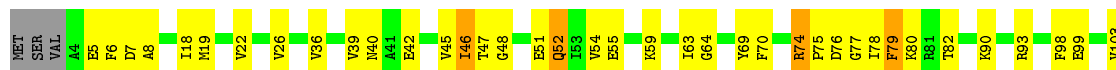
• Molecule 10: 50S ribosomal protein L10e





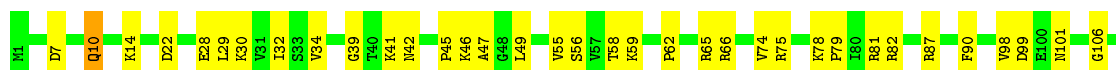
• Molecule 11: 50S ribosomal protein L13P

Chain K: 58% 35% 5%



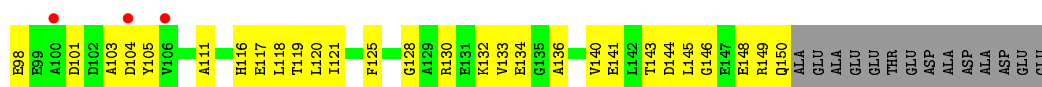
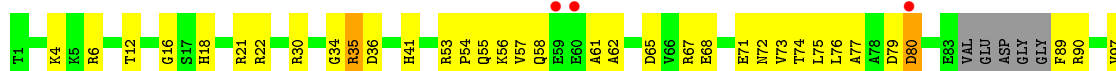
• Molecule 12: 50S ribosomal protein L14P

Chain L: 66% 33%



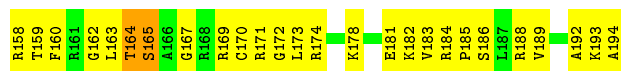
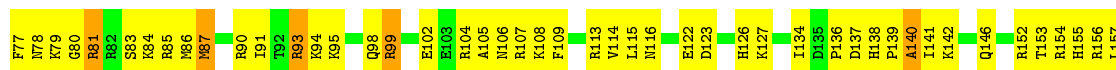
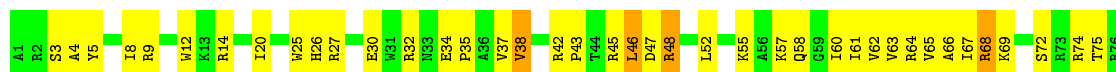
• Molecule 13: 50S ribosomal protein L15P

Chain M: 4% 50% 37% 12%

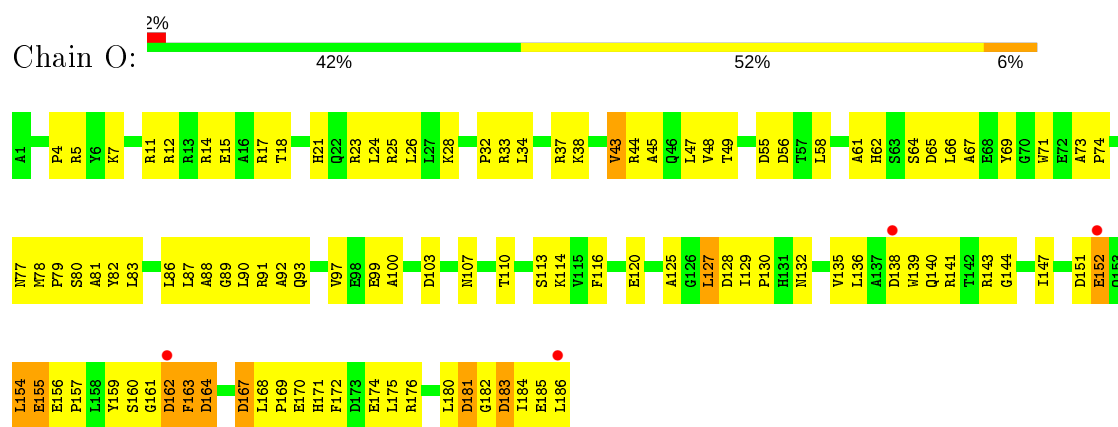


• Molecule 14: 50S ribosomal protein L15E

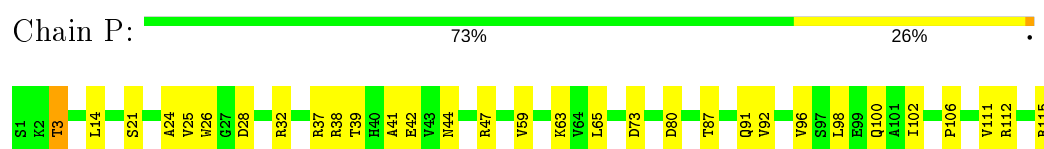
Chain N: 42% 53% 6%



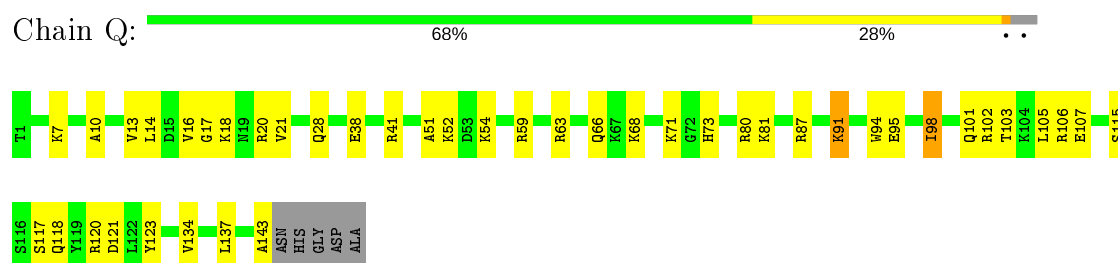
• Molecule 15: 50S ribosomal protein L18P



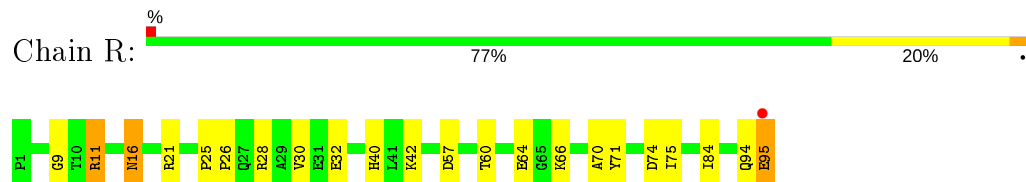
- Molecule 16: 50S ribosomal protein L18E



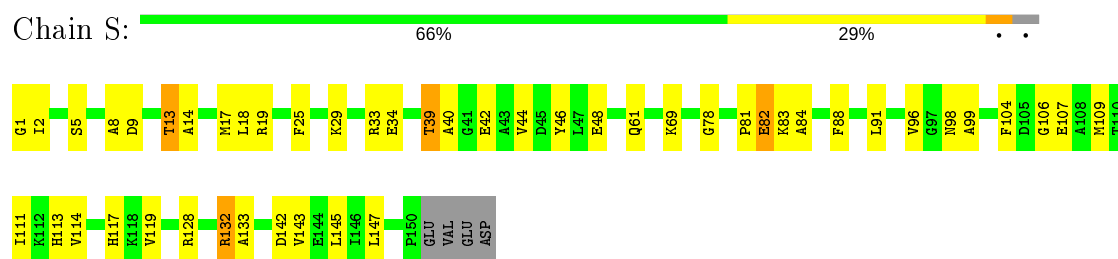
- Molecule 17: 50S ribosomal protein L19E



- Molecule 18: 50S ribosomal protein L21e



- Molecule 19: 50S ribosomal protein L22P



- Molecule 20: 50S ribosomal protein L23P





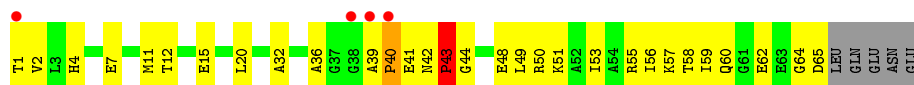
- Molecule 21: 50S ribosomal protein L24P



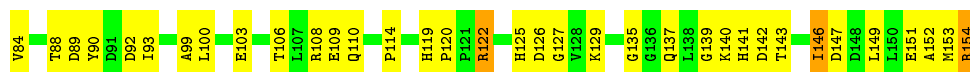
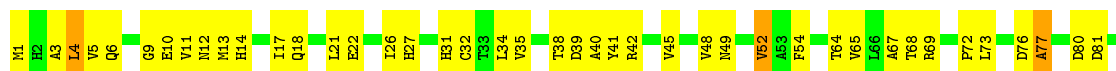
- Molecule 22: 50S ribosomal protein L24E



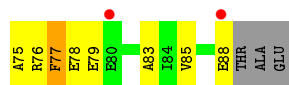
- Molecule 23: 50S ribosomal protein L29P



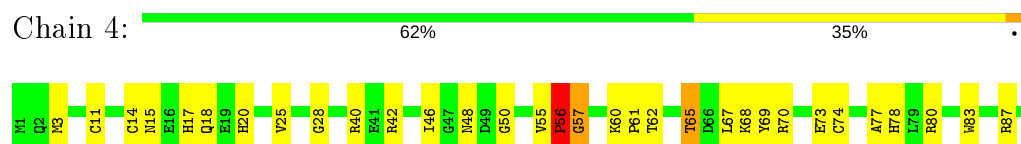
- Molecule 24: 50S ribosomal protein L30P



- Molecule 25: 50S ribosomal protein L31E



- Molecule 26: 50S ribosomal protein L32E



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	211.72Å 299.75Å 573.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.00 49.96 – 2.87	Depositor EDS
% Data completeness (in resolution range)	96.5 (50.00-3.00) 92.7 (49.96-2.87)	Depositor EDS
$R_{merge}$	0.25	Depositor
$R_{sym}$	0.25	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.68 (at 2.86Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.176 , 0.209 0.174 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.1	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 61.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	98536	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CLM, MG, CL, NA, K, CD

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	0.47	5/66076 (0.0%)	0.73	32/103052 (0.0%)
2	B	0.49	0/2905	0.77	4/4528 (0.1%)
3	C	0.44	0/1787	0.73	0/2409
4	D	0.41	0/2689	0.69	0/3652
5	E	0.46	0/1883	0.69	0/2551
6	F	0.35	0/1111	0.62	0/1498
7	G	0.40	0/1382	0.62	0/1880
8	H	0.37	0/896	0.59	0/1219
9	I	0.34	0/241	0.51	0/324
10	J	0.44	0/1246	0.77	1/1686 (0.1%)
11	K	0.44	0/1135	0.64	0/1530
12	L	0.42	0/1003	0.70	0/1351
13	M	0.40	0/1126	0.70	0/1504
14	N	0.47	0/1633	0.74	0/2180
15	O	0.35	0/1473	0.67	0/1999
16	P	0.40	0/873	0.67	0/1181
17	Q	0.41	0/1143	0.58	0/1521
18	R	0.46	0/748	0.73	0/1005
19	S	0.44	0/1172	0.71	0/1578
20	T	0.39	0/648	0.63	1/875 (0.1%)
21	U	0.39	0/957	0.68	0/1289
22	V	0.38	0/417	0.62	0/562
23	W	0.34	0/502	0.58	0/675
24	X	0.42	0/1218	0.69	0/1655
25	Y	0.41	0/664	0.64	0/895
26	Z	0.45	0/1146	0.68	0/1536
27	1	0.46	0/575	0.76	0/763
28	2	0.49	0/437	0.75	0/578
29	3	0.39	0/398	0.57	0/527
30	4	0.47	0/771	0.66	0/1024
All	All	0.46	5/98255 (0.0%)	0.71	38/147027 (0.0%)



Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	90
2	B	1	3
All	All	2	93

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2097	G	O3'-P	8.81	1.71	1.61
1	A	2097	G	C3'-O3'	-7.93	1.31	1.42
1	A	2098	C	P-O5'	-7.79	1.51	1.59
1	A	1206	U	P-OP2	6.71	1.60	1.49
1	A	2098	C	P-OP1	-5.46	1.39	1.49

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1164	U	OP1-P-O3'	-18.95	63.51	105.20
1	A	1164	U	OP2-P-O3'	-17.97	65.67	105.20
1	A	1165	G	O5'-P-OP1	-14.27	92.85	105.70
2	B	3024	U	C2'-C3'-O3'	9.63	130.69	109.50
1	A	1563	G	C2'-C3'-O3'	9.41	130.20	109.50

All (2) chirality outliers are listed below:

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

5 of 93 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	U	Sidechain
1	A	118	G	Sidechain
1	A	22	U	Sidechain
1	A	24	G	Sidechain
1	A	44	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	A	59017	0	29806	952	0
2	B	2600	0	1326	73	0
3	C	1754	0	1763	128	0
4	D	2624	0	2533	176	0
5	E	1858	0	1816	125	0
6	F	1094	0	1085	127	0
7	G	1357	0	1266	70	0
8	H	885	0	854	64	0
9	I	240	0	231	15	0
10	J	1215	0	1215	152	0
11	K	1119	0	1098	68	0
12	L	993	0	1027	50	0
13	M	1114	0	1072	61	0
14	N	1605	0	1676	160	0
15	O	1444	0	1401	117	0
16	P	864	0	873	32	0
17	Q	1133	0	1127	46	0
18	R	734	0	728	18	0
19	S	1149	0	1122	55	0
20	T	641	0	605	20	0
21	U	949	0	923	55	0
22	V	410	0	365	30	0
23	W	499	0	511	30	0
24	X	1195	0	1137	96	0
25	Y	654	0	653	49	0
26	Z	1130	0	1133	62	0
27	1	563	0	598	59	0
28	2	430	0	426	24	0
29	3	393	0	406	25	0
30	4	755	0	728	35	0
31	1	1	0	0	0	0
31	4	1	0	0	0	0
31	A	109	0	0	0	0
31	B	1	0	0	0	0
31	C	1	0	0	0	0
31	D	1	0	0	0	0
31	L	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
31	U	1	0	0	0	0
31	Z	1	0	0	0	0
32	A	2	0	0	0	0
33	A	71	0	0	0	0
33	B	2	0	0	0	0
33	C	1	0	0	0	0
33	E	1	0	0	0	0
33	J	2	0	0	0	0
33	K	1	0	0	0	0
33	M	1	0	0	0	0
33	N	1	0	0	0	0
33	R	1	0	0	0	0
33	S	2	0	0	0	0
33	T	1	0	0	0	0
33	U	1	0	0	0	0
34	4	1	0	0	0	0
34	A	9	0	0	0	0
34	C	1	0	0	0	0
34	D	1	0	0	1	0
34	K	3	0	0	0	0
34	M	1	0	0	0	0
34	N	1	0	0	1	0
34	O	1	0	0	0	0
34	P	1	0	0	0	0
34	R	1	0	0	0	0
34	S	1	0	0	0	0
34	Z	1	0	0	0	0
35	A	20	0	11	0	0
36	1	1	0	0	1	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	37	0	0	9	0
37	2	54	0	0	2	0
37	3	43	0	0	4	0
37	4	68	0	0	11	0
37	A	5906	0	0	192	0
37	B	143	0	0	14	0
37	C	123	0	0	19	0
37	D	147	0	0	32	0
37	E	167	0	0	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
37	F	50	0	0	19	0
37	G	45	0	0	8	0
37	H	29	0	0	7	0
37	I	21	0	0	1	0
37	J	76	0	0	24	0
37	K	54	0	0	3	0
37	L	61	0	0	8	0
37	M	78	0	0	16	0
37	N	125	0	0	19	0
37	O	69	0	0	13	0
37	P	45	0	0	7	0
37	Q	73	0	0	6	0
37	R	53	0	0	3	0
37	S	85	0	0	7	0
37	T	35	0	0	4	0
37	U	40	0	0	7	0
37	V	28	0	0	6	0
37	W	17	0	0	2	0
37	X	69	0	0	17	0
37	Y	28	0	0	6	0
37	Z	98	0	0	18	0
All	All	98536	0	59515	2717	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 2717 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)
10:J:86:ARG:NH1	10:J:133:ILE:HG13	1.57	1.18
5:E:236:THR:HG22	5:E:239:ALA:H	1.03	1.15
1:A:156:C:H5''	14:N:171:ARG:HD3	1.29	1.15
27:1:38:LYS:HE2	27:1:45:LYS:HE2	1.33	1.07
1:A:1160:G:H5'	1:A:1161:A:H5'	1.32	1.06

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%)	210 (89%)	20 (8%)	5 (2%)	7	33
4	D	335/337 (99%)	304 (91%)	24 (7%)	7 (2%)	7	33
5	E	244/246 (99%)	219 (90%)	24 (10%)	1 (0%)	34	72
6	F	134/176 (76%)	96 (72%)	29 (22%)	9 (7%)	1	6
7	G	170/177 (96%)	162 (95%)	8 (5%)	0	100	100
8	H	117/119 (98%)	105 (90%)	10 (8%)	2 (2%)	9	39
9	I	25/348 (7%)	24 (96%)	1 (4%)	0	100	100
10	J	152/167 (91%)	135 (89%)	12 (8%)	5 (3%)	4	21
11	K	140/145 (97%)	129 (92%)	8 (6%)	3 (2%)	7	33
12	L	130/132 (98%)	121 (93%)	7 (5%)	2 (2%)	10	42
13	M	141/164 (86%)	122 (86%)	16 (11%)	3 (2%)	7	33
14	N	192/194 (99%)	170 (88%)	20 (10%)	2 (1%)	15	53
15	O	184/186 (99%)	167 (91%)	10 (5%)	7 (4%)	3	18
16	P	113/115 (98%)	108 (96%)	5 (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%)	138 (93%)	9 (6%)	1 (1%)	22	60
20	T	79/84 (94%)	75 (95%)	4 (5%)	0	100	100
21	U	117/119 (98%)	110 (94%)	6 (5%)	1 (1%)	17	55
22	V	51/66 (77%)	46 (90%)	5 (10%)	0	100	100
23	W	63/70 (90%)	57 (90%)	4 (6%)	2 (3%)	4	22
24	X	152/154 (99%)	146 (96%)	4 (3%)	2 (1%)	12	45
25	Y	80/91 (88%)	70 (88%)	9 (11%)	1 (1%)	12	45
26	Z	140/240 (58%)	136 (97%)	4 (3%)	0	100	100
27	1	71/73 (97%)	63 (89%)	6 (8%)	2 (3%)	5	25

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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%)	85 (94%)	3 (3%)	2 (2%)	6	31
All	All	3633/4235 (86%)	3317 (91%)	259 (7%)	57 (2%)	9	40

5 of 57 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	139	ASP
6	F	20	LYS
6	F	93	LEU
6	F	95	THR
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%)	165 (92%)	14 (8%)	12	42
4	D	282/282 (100%)	266 (94%)	16 (6%)	20	56
5	E	193/193 (100%)	178 (92%)	15 (8%)	12	42
6	F	117/147 (80%)	108 (92%)	9 (8%)	13	42
7	G	152/155 (98%)	148 (97%)	4 (3%)	46	78
8	H	92/92 (100%)	92 (100%)	0	100	100
9	I	27/283 (10%)	27 (100%)	0	100	100
10	J	122/122 (100%)	112 (92%)	10 (8%)	11	39
11	K	118/121 (98%)	109 (92%)	9 (8%)	13	43
12	L	106/106 (100%)	103 (97%)	3 (3%)	43	77
13	M	112/126 (89%)	108 (96%)	4 (4%)	35	70
14	N	166/166 (100%)	157 (95%)	9 (5%)	22	57
15	O	149/149 (100%)	143 (96%)	6 (4%)	31	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
16	P	93/93 (100%)	91 (98%)	2 (2%)	52	81
17	Q	113/116 (97%)	110 (97%)	3 (3%)	44	77
18	R	79/79 (100%)	75 (95%)	4 (5%)	24	60
19	S	117/121 (97%)	113 (97%)	4 (3%)	37	72
20	T	71/73 (97%)	71 (100%)	0	100	100
21	U	105/105 (100%)	102 (97%)	3 (3%)	42	76
22	V	44/52 (85%)	44 (100%)	0	100	100
23	W	51/56 (91%)	50 (98%)	1 (2%)	55	83
24	X	130/130 (100%)	122 (94%)	8 (6%)	18	52
25	Y	66/73 (90%)	62 (94%)	4 (6%)	18	53
26	Z	120/195 (62%)	110 (92%)	10 (8%)	11	39
27	1	56/56 (100%)	53 (95%)	3 (5%)	22	57
28	2	46/46 (100%)	46 (100%)	0	100	100
29	3	42/44 (96%)	41 (98%)	1 (2%)	49	79
30	4	79/79 (100%)	77 (98%)	2 (2%)	47	79
All	All	3027/3441 (88%)	2883 (95%)	144 (5%)	25	62

5 of 144 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
10	J	142	VAL
13	M	30	ARG
26	Z	189	ASN
10	J	166	ASN
11	K	112	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 89 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	116	HIS
17	Q	66	GLN
29	3	16	ASN
14	N	26	HIS
15	O	107	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	2747/2922 (94%)	239 (8%)	38 (1%)
2	B	121/122 (99%)	18 (14%)	6 (4%)
All	All	2868/3044 (94%)	257 (8%)	44 (1%)

5 of 257 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 44 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	1352	A
1	A	1856	C
2	B	3024	U
1	A	1377	C
1	A	1563	G

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 232 ligands modelled in this entry, 231 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	$\# Z  > 2$
35	CLM	A	9001	-	19,20,20	1.41	1 (5%)	23,27,27	1.16	3 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
35	CLM	A	9001	-	-	2/20/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
35	A	9001	CLM	C5-C3	-4.46	1.47	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
35	A	9001	CLM	O4-C4-C3	2.43	116.97	111.09
35	A	9001	CLM	C11-C10-C9	-2.31	116.87	120.08
35	A	9001	CLM	C8-C9-N9	-2.10	117.80	119.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

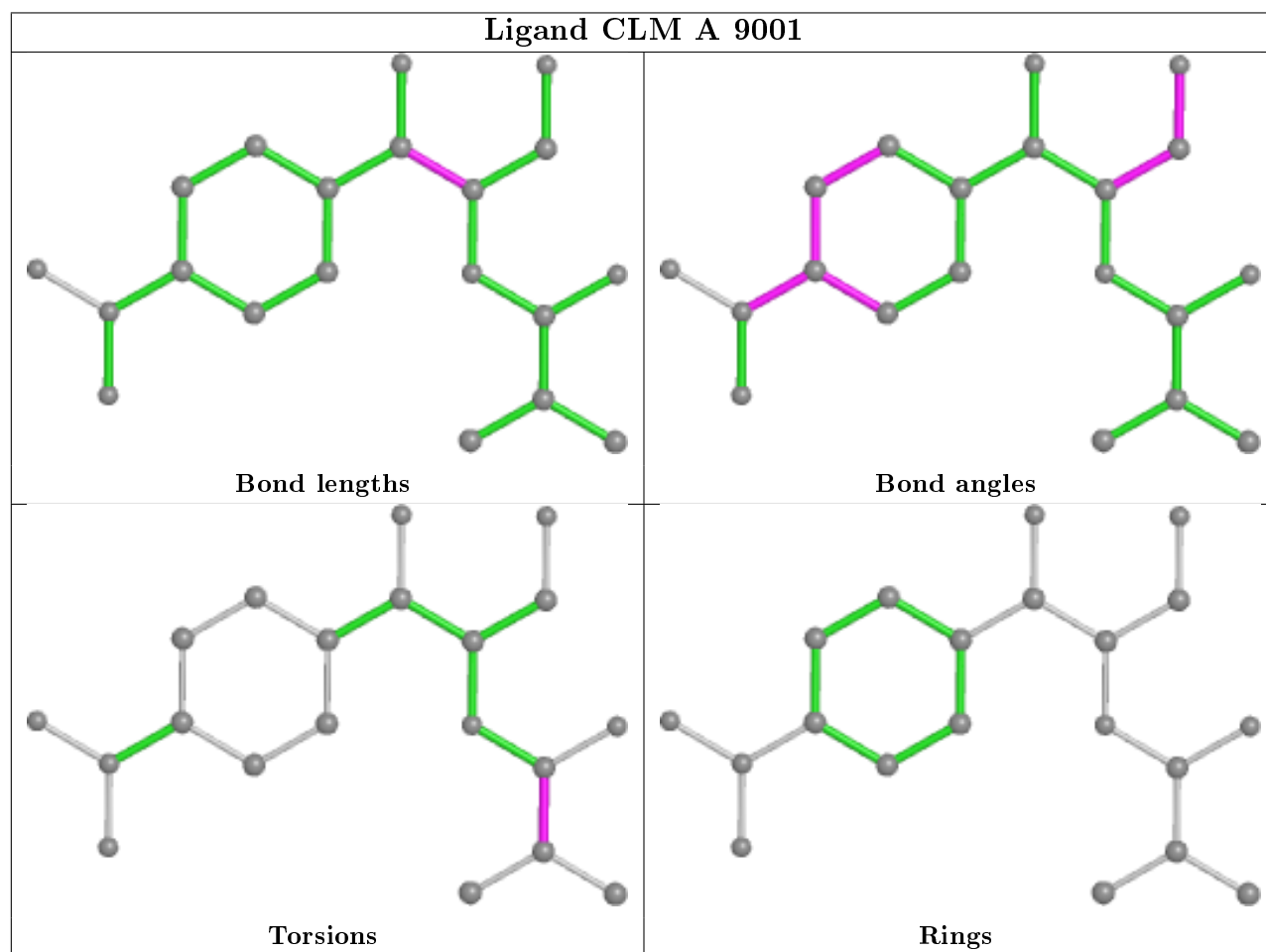
Mol	Chain	Res	Type	Atoms
35	A	9001	CLM	CL2-C1-C2-O2
35	A	9001	CLM	CL2-C1-C2-N2

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	2754/2922 (94%)	-0.34	35 (1%) 77 51	14, 38, 82, 130	0
2	B	122/122 (100%)	-0.09	6 (4%) 29 11	31, 55, 80, 139	0
3	C	237/239 (99%)	-0.17	3 (1%) 77 51	21, 42, 74, 95	0
4	D	337/337 (100%)	-0.30	1 (0%) 94 84	20, 49, 73, 84	0
5	E	246/246 (100%)	-0.42	0 100 100	16, 39, 62, 72	0
6	F	140/176 (79%)	1.22	31 (22%) 0 0	46, 89, 107, 111	0
7	G	172/177 (97%)	0.29	3 (1%) 70 41	39, 61, 80, 85	0
8	H	119/119 (100%)	0.07	2 (1%) 70 41	39, 61, 86, 93	0
9	I	29/348 (8%)	1.50	5 (17%) 1 0	65, 82, 89, 92	0
10	J	156/167 (93%)	-0.07	1 (0%) 89 72	31, 51, 77, 81	0
11	K	142/145 (97%)	-0.25	0 100 100	29, 43, 64, 83	0
12	L	132/132 (100%)	-0.45	0 100 100	28, 45, 63, 70	0
13	M	145/164 (88%)	0.14	6 (4%) 37 14	17, 57, 93, 107	0
14	N	194/194 (100%)	-0.46	0 100 100	22, 36, 55, 65	0
15	O	186/186 (100%)	0.06	4 (2%) 62 33	31, 54, 93, 107	0
16	P	115/115 (100%)	-0.29	0 100 100	31, 47, 63, 67	0
17	Q	143/148 (96%)	-0.13	0 100 100	30, 47, 59, 68	0
18	R	95/95 (100%)	-0.40	1 (1%) 80 56	27, 37, 53, 66	0
19	S	150/154 (97%)	-0.32	0 100 100	24, 39, 58, 66	0
20	T	81/84 (96%)	-0.32	0 100 100	37, 51, 70, 72	0
21	U	119/119 (100%)	0.00	3 (2%) 57 29	33, 49, 73, 86	0
22	V	53/66 (80%)	-0.05	0 100 100	33, 49, 65, 73	0
23	W	65/70 (92%)	0.68	4 (6%) 20 7	43, 64, 99, 103	0
24	X	154/154 (100%)	-0.41	0 100 100	28, 41, 58, 68	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
25	Y	82/91 (90%)	0.20	2 (2%) 59 30	36, 51, 74, 93	0
26	Z	142/240 (59%)	-0.28	3 (2%) 63 34	18, 38, 59, 78	0
27	1	73/73 (100%)	-0.31	0 100 100	36, 52, 68, 75	0
28	2	56/56 (100%)	-0.64	0 100 100	18, 27, 32, 41	0
29	3	46/48 (95%)	0.04	1 (2%) 62 33	28, 52, 78, 91	0
30	4	92/92 (100%)	-0.06	0 100 100	26, 47, 61, 73	0
All	All	6577/7279 (90%)	-0.21	111 (1%) 70 41	14, 44, 83, 139	0

The worst 5 of 111 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
23	W	1	THR	8.4
2	B	3001	U	5.9
2	B	3025	G	5.7
25	Y	88	GLU	4.6
6	F	57	THR	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 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. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
33	NA	A	8384	1/1	0.39	0.56	59,59,59,59	0
33	NA	A	8371	1/1	0.39	0.49	55,55,55,55	0
33	NA	A	8357	1/1	0.57	0.17	56,56,56,56	0
33	NA	A	8364	1/1	0.62	0.27	33,33,33,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
31	MG	A	8049	1/1	0.62	0.51	69,69,69,69	0
33	NA	A	8307	1/1	0.66	0.25	51,51,51,51	0
33	NA	A	8329	1/1	0.67	0.24	48,48,48,48	0
33	NA	A	8382	1/1	0.69	0.54	74,74,74,74	0
34	CL	K	8502	1/1	0.72	0.23	72,72,72,72	0
31	MG	A	8088	1/1	0.74	0.24	44,44,44,44	0
31	MG	A	8089	1/1	0.75	0.13	68,68,68,68	0
33	NA	J	8322	1/1	0.75	0.37	63,63,63,63	0
33	NA	A	8333	1/1	0.75	0.12	24,24,24,24	0
31	MG	A	8022	1/1	0.77	0.11	26,26,26,26	0
35	CLM	A	9001	20/20	0.78	0.31	57,61,68,69	0
33	NA	A	8376	1/1	0.78	0.46	53,53,53,53	0
33	NA	A	8377	1/1	0.78	0.70	66,66,66,66	0
33	NA	A	8373	1/1	0.79	0.30	49,49,49,49	0
33	NA	A	8359	1/1	0.79	0.38	49,49,49,49	0
31	MG	A	8050	1/1	0.79	0.17	79,79,79,79	0
31	MG	A	8114	1/1	0.79	0.13	52,52,52,52	0
31	MG	A	8068	1/1	0.80	0.11	61,61,61,61	0
33	NA	B	8383	1/1	0.80	0.25	45,45,45,45	0
33	NA	A	8326	1/1	0.81	0.27	42,42,42,42	0
33	NA	A	8340	1/1	0.81	0.60	41,41,41,41	0
33	NA	T	8312	1/1	0.81	0.16	44,44,44,44	0
33	NA	S	8386	1/1	0.81	0.50	83,83,83,83	0
33	NA	A	8324	1/1	0.82	0.20	46,46,46,46	0
31	MG	A	8087	1/1	0.82	0.24	71,71,71,71	0
34	CL	4	8504	1/1	0.83	0.23	72,72,72,72	0
33	NA	A	8368	1/1	0.83	0.27	63,63,63,63	0
33	NA	E	8304	1/1	0.83	0.16	30,30,30,30	0
33	NA	A	8363	1/1	0.83	0.33	64,64,64,64	0
33	NA	A	8352	1/1	0.83	0.30	51,51,51,51	0
33	NA	A	8310	1/1	0.84	0.29	21,21,21,21	0
31	MG	A	8113	1/1	0.84	0.19	35,35,35,35	0
31	MG	A	8103	1/1	0.84	0.17	56,56,56,56	0
34	CL	Z	8520	1/1	0.84	0.16	39,39,39,39	0
33	NA	A	8369	1/1	0.85	0.26	51,51,51,51	0
31	MG	A	8076	1/1	0.85	0.18	71,71,71,71	0
34	CL	A	8503	1/1	0.85	0.22	49,49,49,49	0
31	MG	B	8095	1/1	0.85	0.14	61,61,61,61	0
33	NA	A	8342	1/1	0.85	0.24	40,40,40,40	0
33	NA	B	8351	1/1	0.86	0.14	49,49,49,49	0
33	NA	A	8355	1/1	0.86	0.62	57,57,57,57	0
31	MG	U	8073	1/1	0.87	0.09	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
31	MG	A	8072	1/1	0.87	0.18	56,56,56,56	0
31	MG	A	8085	1/1	0.87	0.17	59,59,59,59	0
33	NA	A	8306	1/1	0.87	0.34	51,51,51,51	0
33	NA	A	8316	1/1	0.87	0.28	53,53,53,53	0
33	NA	A	8372	1/1	0.88	0.31	62,62,62,62	0
33	NA	A	8385	1/1	0.88	0.35	45,45,45,45	0
33	NA	A	8315	1/1	0.88	0.22	34,34,34,34	0
31	MG	A	8028	1/1	0.88	0.08	28,28,28,28	0
31	MG	1	8105	1/1	0.88	0.24	28,28,28,28	0
33	NA	A	8302	1/1	0.89	0.28	34,34,34,34	0
33	NA	A	8362	1/1	0.89	0.28	65,65,65,65	0
33	NA	K	8346	1/1	0.89	0.14	43,43,43,43	0
33	NA	A	8378	1/1	0.89	0.86	47,47,47,47	0
33	NA	A	8370	1/1	0.89	0.23	55,55,55,55	0
31	MG	A	8066	1/1	0.89	0.47	38,38,38,38	0
33	NA	A	8375	1/1	0.89	0.24	48,48,48,48	0
31	MG	A	8051	1/1	0.89	0.13	61,61,61,61	0
31	MG	A	8046	1/1	0.89	0.08	45,45,45,45	0
33	NA	C	8345	1/1	0.89	0.15	49,49,49,49	0
34	CL	A	8517	1/1	0.90	0.14	51,51,51,51	0
33	NA	A	8381	1/1	0.90	0.21	45,45,45,45	0
31	MG	A	8042	1/1	0.90	0.22	41,41,41,41	0
33	NA	A	8330	1/1	0.90	0.30	27,27,27,27	0
33	NA	A	8308	1/1	0.90	0.15	49,49,49,49	0
34	CL	D	8519	1/1	0.90	0.30	53,53,53,53	0
31	MG	A	8082	1/1	0.91	0.13	66,66,66,66	0
31	MG	A	8106	1/1	0.91	0.17	51,51,51,51	0
33	NA	A	8319	1/1	0.91	0.11	40,40,40,40	0
31	MG	A	8063	1/1	0.91	0.07	74,74,74,74	0
31	MG	4	8078	1/1	0.91	0.12	35,35,35,35	0
33	NA	A	8327	1/1	0.91	0.17	47,47,47,47	0
33	NA	A	8332	1/1	0.91	0.24	28,28,28,28	0
31	MG	A	8037	1/1	0.91	0.06	39,39,39,39	0
33	NA	S	8337	1/1	0.91	0.11	35,35,35,35	0
31	MG	A	8101	1/1	0.91	0.15	55,55,55,55	0
31	MG	A	8090	1/1	0.92	0.22	46,46,46,46	0
33	NA	M	8380	1/1	0.92	0.55	57,57,57,57	0
33	NA	A	8339	1/1	0.92	0.18	17,17,17,17	0
33	NA	A	8335	1/1	0.92	0.25	53,53,53,53	0
31	MG	A	8108	1/1	0.92	0.07	77,77,77,77	0
32	K	A	8202	1/1	0.92	0.20	56,56,56,56	0
31	MG	A	8045	1/1	0.92	0.11	63,63,63,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	A	8374	1/1	0.92	0.62	56,56,56,56	0
31	MG	A	8102	1/1	0.92	0.18	60,60,60,60	0
33	NA	A	8350	1/1	0.92	0.28	44,44,44,44	0
33	NA	A	8379	1/1	0.93	0.16	51,51,51,51	0
31	MG	A	8081	1/1	0.93	0.07	47,47,47,47	0
31	MG	A	8111	1/1	0.93	0.09	54,54,54,54	0
36	CD	V	8401	1/1	0.93	0.08	63,63,63,63	0
33	NA	A	8341	1/1	0.93	0.15	25,25,25,25	0
34	CL	A	8505	1/1	0.93	0.14	53,53,53,53	0
34	CL	N	8518	1/1	0.93	0.21	41,41,41,41	0
31	MG	A	8024	1/1	0.94	0.10	18,18,18,18	0
33	NA	A	8334	1/1	0.94	0.05	27,27,27,27	0
31	MG	A	8057	1/1	0.94	0.13	38,38,38,38	0
31	MG	A	8100	1/1	0.94	0.12	80,80,80,80	0
31	MG	A	8047	1/1	0.94	0.10	65,65,65,65	0
31	MG	A	8094	1/1	0.94	0.12	61,61,61,61	0
31	MG	A	8044	1/1	0.94	0.15	42,42,42,42	0
31	MG	A	8091	1/1	0.94	0.05	37,37,37,37	0
31	MG	A	8112	1/1	0.94	0.17	49,49,49,49	0
33	NA	A	8313	1/1	0.94	0.22	66,66,66,66	0
34	CL	K	8521	1/1	0.94	0.16	50,50,50,50	0
31	MG	Z	8109	1/1	0.94	0.11	33,33,33,33	0
33	NA	A	8305	1/1	0.94	0.12	32,32,32,32	0
34	CL	A	8515	1/1	0.94	0.26	67,67,67,67	0
33	NA	A	8318	1/1	0.94	0.17	53,53,53,53	0
33	NA	A	8328	1/1	0.94	0.28	30,30,30,30	0
34	CL	C	8509	1/1	0.94	0.19	67,67,67,67	0
31	MG	A	8075	1/1	0.94	0.08	52,52,52,52	0
31	MG	A	8023	1/1	0.94	0.08	36,36,36,36	0
33	NA	U	8343	1/1	0.95	0.10	34,34,34,34	0
33	NA	A	8344	1/1	0.95	0.05	15,15,15,15	0
31	MG	A	8093	1/1	0.95	0.09	40,40,40,40	0
31	MG	D	8055	1/1	0.95	0.07	48,48,48,48	0
31	MG	A	8062	1/1	0.95	0.10	58,58,58,58	0
33	NA	A	8360	1/1	0.95	0.41	42,42,42,42	0
34	CL	A	8514	1/1	0.95	0.12	47,47,47,47	0
31	MG	A	8115	1/1	0.95	0.08	44,44,44,44	0
31	MG	A	8061	1/1	0.95	0.09	35,35,35,35	0
31	MG	A	8015	1/1	0.95	0.05	41,41,41,41	0
31	MG	A	8053	1/1	0.95	0.12	49,49,49,49	0
31	MG	A	8099	1/1	0.95	0.22	48,48,48,48	0
31	MG	A	8011	1/1	0.95	0.07	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
33	NA	A	8354	1/1	0.95	0.14	32,32,32,32	0
33	NA	A	8361	1/1	0.95	0.20	58,58,58,58	0
31	MG	A	8059	1/1	0.95	0.12	37,37,37,37	0
33	NA	A	8353	1/1	0.95	0.12	27,27,27,27	0
31	MG	A	8018	1/1	0.95	0.11	50,50,50,50	0
31	MG	A	8043	1/1	0.95	0.11	41,41,41,41	0
33	NA	A	8349	1/1	0.95	0.22	37,37,37,37	0
33	NA	A	8367	1/1	0.95	0.09	42,42,42,42	0
33	NA	A	8311	1/1	0.95	0.17	51,51,51,51	0
33	NA	A	8317	1/1	0.95	0.12	29,29,29,29	0
34	CL	R	8511	1/1	0.95	0.13	49,49,49,49	0
31	MG	A	8104	1/1	0.96	0.14	42,42,42,42	0
33	NA	A	8338	1/1	0.96	0.14	43,43,43,43	0
33	NA	A	8301	1/1	0.96	0.17	24,24,24,24	0
33	NA	A	8366	1/1	0.96	0.15	53,53,53,53	0
31	MG	C	8065	1/1	0.96	0.11	50,50,50,50	0
31	MG	A	8092	1/1	0.96	0.13	89,89,89,89	0
31	MG	A	8107	1/1	0.96	0.04	34,34,34,34	0
31	MG	A	8110	1/1	0.96	0.05	34,34,34,34	0
34	CL	P	8508	1/1	0.96	0.14	71,71,71,71	0
33	NA	A	8314	1/1	0.96	0.08	28,28,28,28	0
36	CD	4	8404	1/1	0.96	0.07	59,59,59,59	0
31	MG	A	8041	1/1	0.96	0.20	66,66,66,66	0
31	MG	A	8004	1/1	0.96	0.09	33,33,33,33	0
33	NA	A	8365	1/1	0.96	0.21	33,33,33,33	0
33	NA	A	8356	1/1	0.96	0.22	40,40,40,40	0
33	NA	A	8325	1/1	0.96	0.39	58,58,58,58	0
36	CD	1	8403	1/1	0.96	0.10	56,56,56,56	0
31	MG	A	8097	1/1	0.96	0.09	46,46,46,46	0
34	CL	A	8516	1/1	0.96	0.19	69,69,69,69	0
34	CL	O	8507	1/1	0.96	0.19	61,61,61,61	0
31	MG	A	8038	1/1	0.96	0.10	31,31,31,31	0
31	MG	A	8096	1/1	0.96	0.06	42,42,42,42	0
33	NA	J	8309	1/1	0.96	0.07	24,24,24,24	0
31	MG	A	8064	1/1	0.96	0.17	20,20,20,20	0
31	MG	A	8033	1/1	0.96	0.05	31,31,31,31	0
34	CL	A	8513	1/1	0.96	0.11	52,52,52,52	0
33	NA	R	8348	1/1	0.96	0.08	27,27,27,27	0
31	MG	A	8077	1/1	0.97	0.09	40,40,40,40	0
33	NA	A	8303	1/1	0.97	0.17	41,41,41,41	0
31	MG	A	8001	1/1	0.97	0.04	22,22,22,22	0
31	MG	A	8040	1/1	0.97	0.11	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
31	MG	A	8067	1/1	0.97	0.13	49,49,49,49	0
31	MG	L	8069	1/1	0.97	0.08	64,64,64,64	0
34	CL	K	8501	1/1	0.97	0.18	61,61,61,61	0
33	NA	A	8321	1/1	0.97	0.23	45,45,45,45	0
31	MG	A	8058	1/1	0.97	0.12	43,43,43,43	0
34	CL	M	8510	1/1	0.97	0.12	60,60,60,60	0
31	MG	A	8009	1/1	0.97	0.06	32,32,32,32	0
31	MG	A	8013	1/1	0.97	0.12	48,48,48,48	0
32	K	A	8201	1/1	0.97	0.14	55,55,55,55	0
34	CL	A	8522	1/1	0.97	0.18	66,66,66,66	0
36	CD	P	8405	1/1	0.97	0.04	90,90,90,90	0
33	NA	A	8323	1/1	0.97	0.14	37,37,37,37	0
31	MG	A	8071	1/1	0.97	0.08	72,72,72,72	0
31	MG	A	8052	1/1	0.97	0.18	52,52,52,52	0
31	MG	A	8054	1/1	0.97	0.11	38,38,38,38	0
31	MG	A	8117	1/1	0.97	0.09	25,25,25,25	0
31	MG	A	8005	1/1	0.97	0.10	36,36,36,36	0
31	MG	A	8080	1/1	0.97	0.08	42,42,42,42	0
31	MG	A	8017	1/1	0.97	0.04	22,22,22,22	0
31	MG	A	8021	1/1	0.97	0.09	36,36,36,36	0
31	MG	A	8098	1/1	0.97	0.13	29,29,29,29	0
31	MG	A	8020	1/1	0.97	0.08	36,36,36,36	0
33	NA	N	8347	1/1	0.97	0.08	17,17,17,17	0
31	MG	A	8070	1/1	0.97	0.09	47,47,47,47	0
31	MG	A	8035	1/1	0.97	0.04	39,39,39,39	0
31	MG	A	8048	1/1	0.97	0.06	43,43,43,43	0
31	MG	A	8032	1/1	0.98	0.05	26,26,26,26	0
31	MG	A	8019	1/1	0.98	0.05	27,27,27,27	0
31	MG	A	8031	1/1	0.98	0.03	22,22,22,22	0
31	MG	A	8036	1/1	0.98	0.03	39,39,39,39	0
34	CL	S	8506	1/1	0.98	0.12	43,43,43,43	0
31	MG	A	8007	1/1	0.98	0.07	26,26,26,26	0
36	CD	2	8402	1/1	0.98	0.07	57,57,57,57	0
33	NA	A	8331	1/1	0.98	0.14	38,38,38,38	0
31	MG	A	8014	1/1	0.98	0.06	30,30,30,30	0
31	MG	A	8016	1/1	0.98	0.10	27,27,27,27	0
31	MG	A	8116	1/1	0.98	0.10	60,60,60,60	0
31	MG	A	8026	1/1	0.98	0.09	27,27,27,27	0
31	MG	A	8083	1/1	0.98	0.07	42,42,42,42	0
31	MG	A	8074	1/1	0.98	0.06	28,28,28,28	0
33	NA	A	8336	1/1	0.98	0.04	34,34,34,34	0
31	MG	A	8030	1/1	0.98	0.06	30,30,30,30	0

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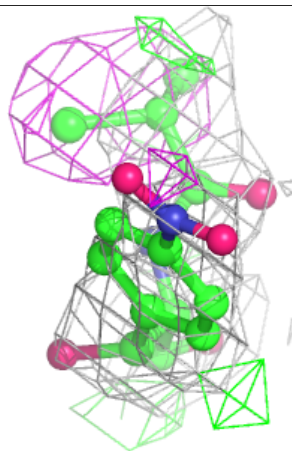
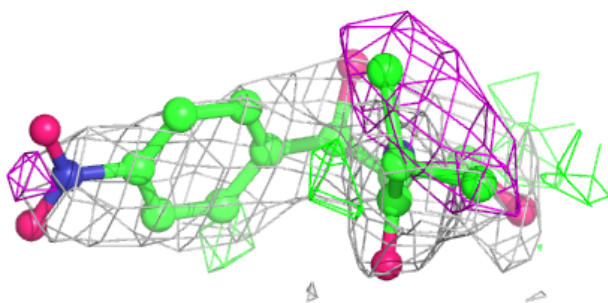
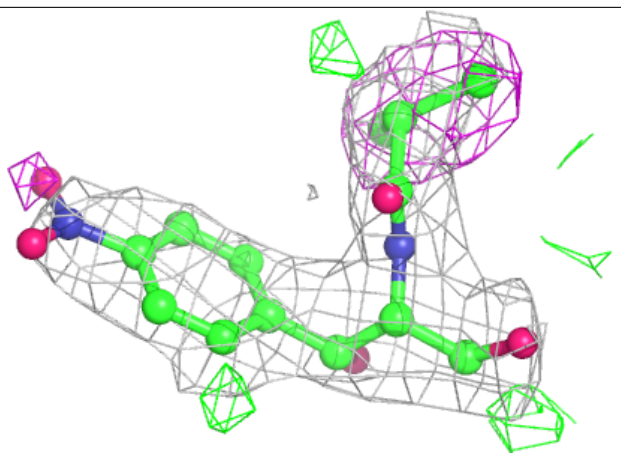
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
31	MG	A	8027	1/1	0.98	0.04	30,30,30,30	0
31	MG	A	8008	1/1	0.98	0.06	43,43,43,43	0
31	MG	A	8034	1/1	0.98	0.04	26,26,26,26	0
31	MG	A	8006	1/1	0.98	0.04	28,28,28,28	0
31	MG	A	8086	1/1	0.98	0.18	35,35,35,35	0
31	MG	A	8029	1/1	0.98	0.07	40,40,40,40	0
31	MG	A	8056	1/1	0.98	0.07	44,44,44,44	0
31	MG	A	8025	1/1	0.98	0.09	47,47,47,47	0
33	NA	A	8320	1/1	0.98	0.07	20,20,20,20	0
31	MG	A	8010	1/1	0.98	0.05	29,29,29,29	0
31	MG	A	8003	1/1	0.99	0.06	18,18,18,18	0
31	MG	A	8084	1/1	0.99	0.10	60,60,60,60	0
31	MG	A	8012	1/1	0.99	0.12	37,37,37,37	0
34	CL	A	8512	1/1	0.99	0.07	41,41,41,41	0
31	MG	A	8079	1/1	0.99	0.09	36,36,36,36	0
31	MG	A	8060	1/1	0.99	0.18	36,36,36,36	0
31	MG	A	8002	1/1	0.99	0.06	38,38,38,38	0
31	MG	A	8039	1/1	0.99	0.05	52,52,52,52	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 CLM A 9001:**

$2mF_o-DF_c$  (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.