



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 11, 2017 – 10:58 PM EDT

PDB ID : 4WOI
Title : 4,5-linked aminoglycoside antibiotics regulate the bacterial ribosome by targeting dynamic conformational processes within intersubunit bridge B2
Authors : Pulk, A.; Cate, J.H.D.; Blanchard, S.; Wasserman, M.; Altman, R.; Zhou, Z.; Zinder, J.; Green, K.; Garneau-Tsodikova, S.
Deposited on : unknown
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

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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20029824

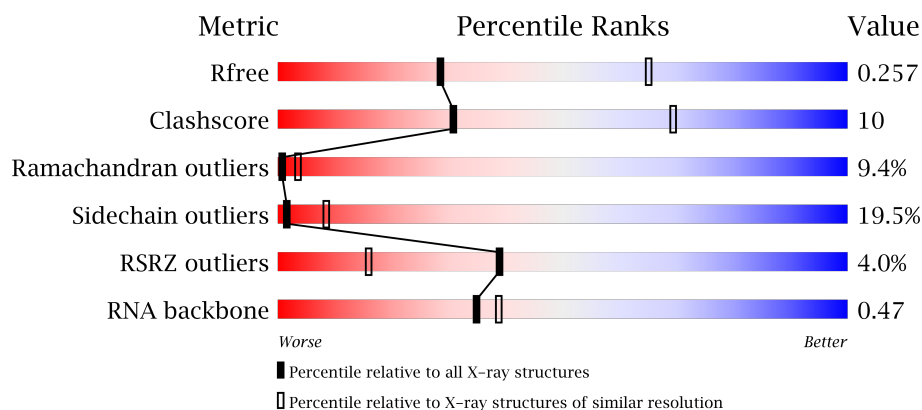
1 Overall quality at a glance

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.






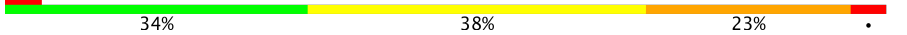

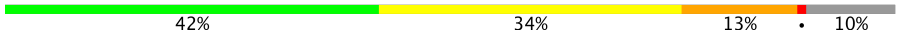
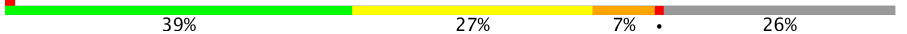
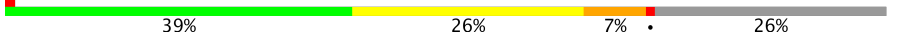
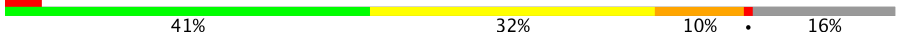



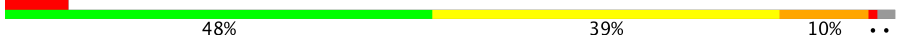
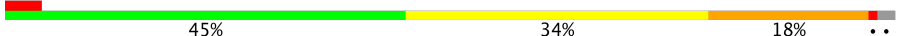











Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)
RNA backbone	2435	1007 (3.34-2.66)

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	AA	1542	<div> <div>54%</div> <div>36%</div> <div>9%</div> <div>.</div> </div>
1	DA	1542	<div> <div>56%</div> <div>34%</div> <div>10%</div> </div>
2	AB	241	<div> <div>12%</div> <div>51%</div> <div>25%</div> <div>12%</div> <div>.</div> <div>10%</div> </div>
2	DB	241	<div> <div>6%</div> <div>36%</div> <div>39%</div> <div>13%</div> <div>.</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
3	AC	233	
3	DC	233	
4	AD	206	
4	DD	206	
5	AE	167	
5	DE	167	
6	AF	135	
6	DF	135	
7	AG	179	
7	DG	179	
8	AH	130	
8	DH	130	
9	AI	130	
9	DI	130	
10	AJ	103	
10	DJ	103	
11	AK	129	
11	DK	129	
12	AL	124	
12	DL	124	
13	AM	118	
13	DM	118	
14	AN	101	
14	DN	101	
15	AO	89	








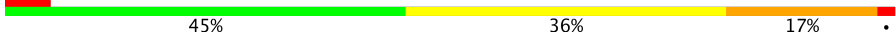

















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Mol	Chain	Length	Quality of chain
15	DO	89	
16	AP	82	
16	DP	82	
17	AQ	84	
17	DQ	84	
18	AR	75	
18	DR	75	
19	AS	92	
19	DS	92	
20	AT	87	
20	DT	87	
21	AU	71	
21	DU	71	
22	AV	185	
23	AW	16	
23	DV	16	
24	AX	76	
24	DW	76	
25	BA	2904	
25	CA	2904	
26	BB	120	
26	CB	120	
27	BC	273	
27	CC	273	
28	BD	209	











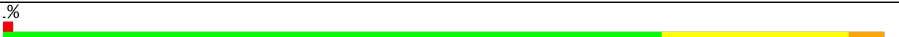


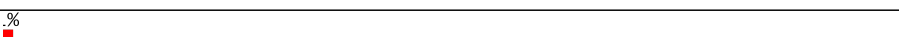
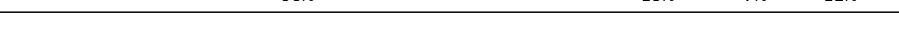
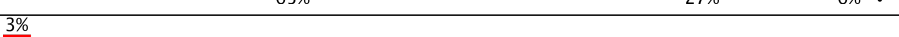



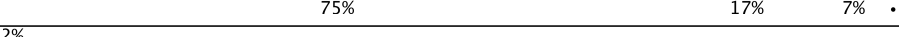





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Mol	Chain	Length	Quality of chain
28	CD	209	
29	BE	201	
29	CE	201	
30	BF	179	
30	CF	179	
31	BG	177	
31	CG	177	
32	BH	149	
32	CH	149	
33	BI	142	
33	CI	142	
34	BJ	142	
34	CJ	142	
35	BK	123	
35	CK	123	
36	BL	144	
36	CL	144	
37	BM	136	
37	CM	136	
38	BN	127	
38	CN	127	
39	BO	117	
39	CO	117	
40	BP	115	
40	CP	115	

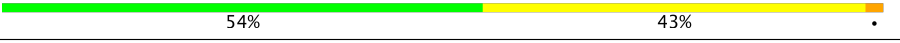



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Mol	Chain	Length	Quality of chain
41	BQ	118	
41	CQ	118	
42	BR	103	
42	CR	103	
43	BS	110	
43	CS	110	
44	BT	100	
44	CT	100	
45	BU	104	
45	CU	104	
46	BV	94	
46	CV	94	
47	BW	85	
47	CW	85	
48	BX	78	
48	CX	78	
49	BY	63	
49	CY	63	
50	BZ	59	
50	CZ	59	
51	B0	57	
51	C0	57	
52	B1	55	
52	C1	55	
53	B2	46	

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Mol	Chain	Length	Quality of chain
53	C2	46	
54	B3	65	
54	C3	65	
55	B4	38	
55	C4	38	
56	B5	228	

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
57	MG	AA	1603	-	-	-	X
57	MG	AA	1612	-	-	-	X
57	MG	AA	1614	-	-	-	X
57	MG	AA	1619	-	-	-	X
57	MG	AA	1631	-	-	-	X
57	MG	AA	1636	-	-	-	X
57	MG	AA	1640	-	-	-	X
57	MG	AA	1645	-	-	-	X
57	MG	AA	1669	-	-	-	X
57	MG	BA	3014	-	-	-	X
57	MG	BA	3021	-	-	-	X
57	MG	BA	3025	-	-	-	X
57	MG	BA	3038	-	-	-	X
57	MG	BA	3044	-	-	-	X
57	MG	BA	3045	-	-	-	X
57	MG	BA	3066	-	-	-	X
57	MG	BA	3067	-	-	-	X
57	MG	BA	3077	-	-	-	X
57	MG	BA	3105	-	-	-	X
57	MG	BA	3109	-	-	-	X
57	MG	BA	3112	-	-	-	X
57	MG	BA	3118	-	-	-	X
57	MG	BA	3120	-	-	-	X
57	MG	BA	3128	-	-	-	X
57	MG	BA	3134	-	-	-	X
57	MG	BA	3138	-	-	-	X
57	MG	BA	3154	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
57	MG	BA	3155	-	-	-	X
57	MG	BA	3157	-	-	-	X
57	MG	BA	3161	-	-	-	X
57	MG	BA	3166	-	-	-	X
57	MG	BA	3176	-	-	-	X
57	MG	BA	3179	-	-	-	X
57	MG	BA	3188	-	-	-	X
57	MG	CA	3013	-	-	-	X
57	MG	CA	3019	-	-	-	X
57	MG	CA	3041	-	-	-	X
57	MG	CA	3048	-	-	-	X
57	MG	CA	3064	-	-	-	X
57	MG	CA	3094	-	-	-	X
57	MG	CA	3098	-	-	-	X
57	MG	CA	3104	-	-	-	X
57	MG	CA	3108	-	-	-	X
57	MG	CA	3109	-	-	-	X
57	MG	CA	3116	-	-	-	X
57	MG	CA	3133	-	-	-	X
57	MG	CA	3142	-	-	-	X
57	MG	CA	3143	-	-	-	X
57	MG	CA	3145	-	-	-	X
57	MG	CA	3151	-	-	-	X
57	MG	CA	3161	-	-	-	X
57	MG	CD	301	-	-	-	X
57	MG	DA	1605	-	-	-	X
57	MG	DA	1633	-	-	-	X
57	MG	DA	1638	-	-	-	X
57	MG	DA	1644	-	-	-	X
58	PAR	AA	1672	-	-	-	X
58	PAR	BA	3001	-	-	-	X
58	PAR	BA	3002	-	-	-	X
58	PAR	BA	3003	-	-	-	X
58	PAR	BA	3004	-	-	-	X
58	PAR	BA	3005	-	-	X	X
58	PAR	CA	3166	-	-	-	X
58	PAR	CA	3167	-	-	-	X
58	PAR	CA	3168	-	-	-	X
58	PAR	CA	3169	-	-	-	X
58	PAR	DA	1654	-	-	-	X
58	PAR	DA	1655	-	-	-	X
59	ZN	B4	9501	-	-	-	X

2 Entry composition

There are 60 unique types of molecules in this entry. The entry contains 294484 atoms, of which 450 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 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	1538	Total	C	N	O	P	0	0	0
			32995	14716	6050	10691	1538			
1	DA	1539	Total	C	N	O	P	0	0	0
			33015	14725	6052	10699	1539			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AB	218	Total	C	N	O	S	0	0	0
			1704	1081	305	311	7			
2	DB	218	Total	C	N	O	S	0	0	0
			1704	1081	305	311	7			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AC	206	Total	C	N	O	S	0	0	0
			1624	1028	305	288	3			
3	DC	206	Total	C	N	O	S	0	0	0
			1624	1028	305	288	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AD	205	Total	C	N	O	S	0	0	0
			1643	1026	315	298	4			
4	DD	205	Total	C	N	O	S	0	0	0
			1643	1026	315	298	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AE	150	Total	C	N	O	S	0	0	0
			1105	687	211	201	6			
5	DE	150	Total	C	N	O	S	0	0	0
			1105	687	211	201	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AF	100	Total	C	N	O	S	0	0	0
			817	515	148	148	6			
6	DF	100	Total	C	N	O	S	0	0	0
			817	515	148	148	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AG	151	Total	C	N	O	S	0	0	0
			1181	735	227	215	4			
7	DG	151	Total	C	N	O	S	0	0	0
			1181	735	227	215	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AH	129	Total	C	N	O	S	0	0	0
			979	616	173	184	6			
8	DH	129	Total	C	N	O	S	0	0	0
			979	616	173	184	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	AI	127	Total	C	N	O	S	0	0	0
			1022	634	206	179	3			
9	DI	127	Total	C	N	O	S	0	0	0
			1022	634	206	179	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AJ	98	Total	C	N	O	S	0	0	0
			786	493	150	142	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	DJ	98	Total	C	N	O	S	0	0	0
			786	493	150	142	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AK	117	Total	C	N	O	S	0	0	0
			877	540	174	160	3			
11	DK	117	Total	C	N	O	S	0	0	0
			877	540	174	160	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AL	123	Total	C	N	O	S	0	0	0
			955	590	196	165	4			
12	DL	123	Total	C	N	O	S	0	0	0
			955	590	196	165	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AM	114	Total	C	N	O	S	0	0	0
			883	546	178	156	3			
13	DM	114	Total	C	N	O	S	0	0	0
			883	546	178	156	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AN	96	Total	C	N	O	S	0	0	0
			774	483	160	128	3			
14	DN	96	Total	C	N	O	S	0	0	0
			774	483	160	128	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AO	88	Total	C	N	O	S	0	0	0
			716	440	146	129	1			
15	DO	88	Total	C	N	O	S	0	0	0
			716	440	146	129	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AP	82	Total	C	N	O	S	0	0	0
			649	406	128	114	1			
16	DP	82	Total	C	N	O	S	0	0	0
			649	406	128	114	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AQ	80	Total	C	N	O	S	0	0	0
			648	411	121	113	3			
17	DQ	80	Total	C	N	O	S	0	0	0
			648	411	121	113	3			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	AR	55	Total	C	N	O	0	0	0
			455	288	86	81			
18	DR	55	Total	C	N	O	0	0	0
			455	288	86	81			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AS	79	Total	C	N	O	S	0	0	0
			637	408	120	107	2			
19	DS	79	Total	C	N	O	S	0	0	0
			637	408	120	107	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AT	85	Total	C	N	O	S	0	0	0
			665	411	137	114	3			
20	DT	85	Total	C	N	O	S	0	0	0
			665	411	137	114	3			

- Molecule 21 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	AU	51	Total	C	N	O	S	0	0	0
			425	265	86	73	1			
21	DU	51	Total	C	N	O	S	0	0	0
			425	265	86	73	1			

- Molecule 22 is a protein called Ribosome-recycling factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	AV	183	Total	C	N	O	S	0	0	0
			1419	871	260	283	5			

- Molecule 23 is a RNA chain called Messenger RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	AW	15	Total	C	N	O	P	0	0	0
			324	145	61	103	15			
23	DV	16	Total	C	N	O	P	0	0	0
			346	155	66	109	16			

- Molecule 24 is a RNA chain called Phenylalanine specific transfer RNA, tRNA-Phe.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	AX	76	Total	C	N	O	P	0	0	0
			1623	723	290	534	76			
24	DW	76	Total	C	N	O	P	0	0	0
			1623	723	290	534	76			

- Molecule 25 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	BA	2897	Total	C	N	O	P	0	0	0
			62195	27745	11446	20107	2897			
25	CA	2897	Total	C	N	O	P	0	0	0
			62195	27745	11446	20107	2897			

- Molecule 26 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	BB	119	Total	C	N	O	P	0	0	0
			2549	1135	466	829	119			
26	CB	118	Total	C	N	O	P	0	0	0
			2529	1126	464	821	118			

- Molecule 27 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	BC	271	Total	C	N	O	S	0	0	0
			2082	1288	423	364	7			
27	CC	271	Total	C	N	O	S	0	0	0
			2082	1288	423	364	7			

- Molecule 28 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	BD	209	Total	C	N	O	S	0	0	0
			1565	979	288	294	4			
28	CD	209	Total	C	N	O	S	0	0	0
			1565	979	288	294	4			

- Molecule 29 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	BE	201	Total	C	N	O	S	0	0	0
			1552	974	283	290	5			
29	CE	201	Total	C	N	O	S	0	0	0
			1552	974	283	290	5			

- Molecule 30 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	BF	177	Total	C	N	O	S	0	0	0
			1410	899	249	256	6			
30	CF	177	Total	C	N	O	S	0	0	0
			1410	899	249	256	6			

- Molecule 31 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	BG	176	Total	C	N	O	S	0	0	0
			1323	832	243	246	2			
31	CG	176	Total	C	N	O	S	0	0	0
			1323	832	243	246	2			

- Molecule 32 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	BH	149	Total	C	N	O	S	0	0	0
			1107	696	197	213	1			
32	CH	149	Total	C	N	O	S	0	0	0
			1107	696	197	213	1			

- Molecule 33 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	BI	141	Total	C	N	O	S	0	0	0
			1032	651	179	196	6			
33	CI	141	Total	C	N	O	S	0	0	0
			1032	651	179	196	6			

- Molecule 34 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	BJ	142	Total	C	N	O	S	0	0	0
			1129	714	212	199	4			
34	CJ	142	Total	C	N	O	S	0	0	0
			1129	714	212	199	4			

- Molecule 35 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BK	122	Total	C	N	O	S	0	0	0
			938	587	180	165	6			
35	CK	122	Total	C	N	O	S	0	0	0
			938	587	180	165	6			

- Molecule 36 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	BL	143	Total	C	N	O	S	0	0	0
			1045	649	206	189	1			
36	CL	143	Total	C	N	O	S	0	0	0
			1045	649	206	189	1			

- Molecule 37 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	BM	136	Total	C	N	O	S	0	0	0
			1074	686	205	177	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	CM	136	Total	C	N	O	S	0	0	0
			1074	686	205	177	6			

- Molecule 38 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	BN	120	Total	C	N	O	S	0	0	0
			960	593	196	166	5			
38	CN	120	Total	C	N	O	S	0	0	0
			960	593	196	166	5			

- Molecule 39 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	BO	116	Total	C	N	O	S	0	0	0
			892	552	178	162				
39	CO	116	Total	C	N	O	S	0	0	0
			892	552	178	162				

- Molecule 40 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	BP	114	Total	C	N	O	S	0	0	0
			917	574	179	163	1			
40	CP	114	Total	C	N	O	S	0	0	0
			917	574	179	163	1			

- Molecule 41 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BQ	117	Total	C	N	O	S	0	0	0
			947	604	192	151				
41	CQ	117	Total	C	N	O	S	0	0	0
			947	604	192	151				

- Molecule 42 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BR	103	Total	C	N	O	S	0	0	0
			816	516	153	145	2			
42	CR	103	Total	C	N	O	S	0	0	0
			816	516	153	145	2			

- Molecule 43 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	BS	110	Total	C	N	O	S	0	0	0
			857	532	166	156	3			
43	CS	110	Total	C	N	O	S	0	0	0
			857	532	166	156	3			

- Molecule 44 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	BT	93	Total	C	N	O	S	0	0	0
			738	466	139	131	2			
44	CT	93	Total	C	N	O	S	0	0	0
			738	466	139	131	2			

- Molecule 45 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
45	BU	102	Total	C	N	O	0	0	0
			779	492	146	141			
45	CU	102	Total	C	N	O	0	0	0
			779	492	146	141			

- Molecule 46 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	BV	94	Total	C	N	O	S	0	0	0
			753	479	137	134	3			
46	CV	94	Total	C	N	O	S	0	0	0
			753	479	137	134	3			

- Molecule 47 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	BW	76	Total	C	N	O	S	0	0	0
			580	359	117	103	1			
47	CW	75	Total	C	N	O	S	0	0	0
			569	353	113	102	1			

- Molecule 48 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	BX	77	Total	C	N	O	S	0	0	0
			625	388	129	106	2			
48	CX	77	Total	C	N	O	S	0	0	0
			625	388	129	106	2			

- Molecule 49 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	BY	63	Total	C	N	O	S	0	0	0
			509	313	99	95	2			
49	CY	63	Total	C	N	O	S	0	0	0
			509	313	99	95	2			

- Molecule 50 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	BZ	58	Total	C	N	O	S	0	0	0
			449	281	87	79	2			
50	CZ	58	Total	C	N	O	S	0	0	0
			449	281	87	79	2			

- Molecule 51 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	B0	56	Total	C	N	O	S	0	0	0
			444	269	94	80	1			
51	C0	56	Total	C	N	O	S	0	0	0
			444	269	94	80	1			

- Molecule 52 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
52	B1	50	Total	C	N	O	0	0	0
			409	263	75	71			
52	C1	50	Total	C	N	O	0	0	0
			409	263	75	71			

- Molecule 53 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	B2	46	Total	C	N	O	S	0	0	0
			377	228	90	57	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	C2	46	Total	C	N	O	S	0	0	0
			377	228	90	57	2			

- Molecule 54 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	B3	64	Total	C	N	O	S	0	0	0
			504	323	105	74	2			
54	C3	64	Total	C	N	O	S	0	0	0
			504	323	105	74	2			

- Molecule 55 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	B4	38	Total	C	N	O	S	0	0	0
			302	185	65	48	4			
55	C4	38	Total	C	N	O	S	0	0	0
			302	185	65	48	4			

- Molecule 56 is a protein called 50S ribosomal protein L1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
56	B5	191	Total	C	N	O	0	0	1
			1142	691	221	230			

- Molecule 57 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
57	BB	4	Total	Mg	0	0
			4	4		
57	BO	1	Total	Mg	0	0
			1	1		
57	BA	189	Total	Mg	0	0
			189	189		
57	CA	165	Total	Mg	0	0
			165	165		
57	CB	3	Total	Mg	0	0
			3	3		
57	BL	2	Total	Mg	0	0
			2	2		
57	DA	53	Total	Mg	0	0
			53	53		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
57	CD	1	Total 1 Mg 1	0	0
57	AA	71	Total 71 Mg 71	0	0
57	BQ	1	Total 1 Mg 1	0	0
57	CQ	1	Total 1 Mg 1	0	0
57	AN	1	Total 1 Mg 1	0	0
57	DN	1	Total 1 Mg 1	0	0
57	DD	2	Total 2 Mg 2	0	0

-
- The chemical structure of PAR (Pantetheine) is a complex molecule consisting of several fused and linked rings. It features a central pyrimidine ring (C22-C27) with an amino group (NH₂, N22) at C27. This ring is linked via an oxygen atom (O11) to a five-membered ring containing a nitrogen atom (N21) and a hydroxyl group (OH, O18). Another oxygen atom (O12) links the central ring to a six-membered ring containing a hydroxyl group (OH, O16) and a nitrogen atom (N23). A third oxygen atom (O13) links the central ring to a five-membered ring containing a hydroxyl group (OH, O23) and a nitrogen atom (N24). The structure is highly branched and contains multiple stereocenters, with atoms labeled with IDs such as C10, C11, C12, C13, C14, C15, C16, C17, C18, C19, C20, C21, C22, C23, C24, C25, C26, C27, C28, C29, C30, C31, C32, C33, C34, C35, C36, C37, C38, C39, C40, C41, C42, C43, C44, C45, C46, C47, C48, C49, C50, C51, C52, C53, C54, C55, C56, C57, C58, C59, C60, C61, C62, C63, C64, C65, C66, C67, C68, C69, C70, C71, C72, C73, C74, C75, C76, C77, C78, C79, C80, C81, C82, C83, C84, C85, C86, C87, C88, C89, C90, C91, C92, C93, C94, C95, C96, C97, C98, C99, C100, C101, C102, C103, C104, C105, C106, C107, C108, C109, C110, C111, C112, C113, C114, C115, C116, C117, C118, C119, C120, C121, C122, C123, C124, C125, C126, C127, C128, C129, C130, C131, C132, C133, C134, C135, C136, C137, C138, C139, C140, C141, C142, C143, C144, C145, C146, C147, C148, C149, C150, C151, C152, C153, C154, C155, C156, C157, C158, C159, C160, C161, C162, C163, C164, C165, C166, C167, C168, C169, C170, C171, C172, C173, C174, C175, C176, C177, C178, C179, C180, C181, C182, C183, C184, C185, C186, C187, C188, C189, C190, C191, C192, C193, C194, C195, C196, C197, C198, C199, C200, C201, C202, C203, C204, C205, C206, C207, C208, C209, C210, C211, C212, C213, C214, C215, C216, C217, C218, C219, C220, C221, C222, C223, C224, C225, C226, C227, C228, C229, C230, C231, C232, C233, C234, C235, C236, C237, C238, C239, C240, C241, C242, C243, C244, C245, C246, C247, C248, C249, C250, C251, C252, C253, C254, C255, C256, C257, C258, C259, C260, C261, C262, C263, C264, C265, C266, C267, C268, C269, C270, C271, C272, C273, C274, C275, C276, C277, C278, C279, C280, C281, C282, C283, C284, C285, C286, C287, C288, C289, C290, C291, C292, C293, C294, C295, C296, C297, C298, C299, C300, C301, C302, C303, C304, C305, C306, C307, C308, C309, C310, C311, C312, C313, C314, C315, C316, C317, C318, C319, C320, C321, C322, C323, C324, C325, C326, C327, C328, C329, C330, C331, C332, C333, C334, C335, C336, C337, C338, C339, C340, C341, C342, C343, C344, C345, C346, C347, C348, C349, C350, C351, C352, C353, C354, C355, C356, C357, C358, C359, C360, C361, C362, C363, C364, C365, C366, C367, C368, C369, C370, C371, C372, C373, C374, C375, C376, C377, C378, C379, C380, C381, C382, C383, C384, C385, C386, C387, C388, C389, C390, C391, C392, C393, C394, C395, C396, C397, C398, C399, C400, C401, C402, C403, C404, C405, C406, C407, C408, C409, C410, C411, C412, C413, C414, C415, C416, C417, C418, C419, C420, C421, C422, C423, C424, C425, C426, C427, C428, C429, C430, C431, C432, C433, C434, C435, C436, C437, C438, C439, C440, C441, C442, C443, C444, C445, C446, C447, C448, C449, C450, C451, C452, C453, C454, C455, C456, C457, C458, C459, C460, C461, C462, C463, C464, C465, C466, C467, C468, C469, C470, C471, C472, C473, C474, C475, C476, C477, C478, C479, C480, C481, C482, C483, C484, C485, C486, C487, C488, C489, C490, C491, C492, C493, C494, C495, C496, C497, C498, C499, C500, C501, C502, C503, C504, C505, C506, C507, C508, C509, C510, C511, C512, C513, C514, C515, C516, C517, C518, C519, C520, C521, C522, C523, C524, C525, C526, C527, C528, C529, C530, C531, C532, C533, C534, C535, C536, C537, C538, C539, C540, C541, C542, C543, C544, C545, C546, C547, C548, C549, C550, C551, C552, C553, C554, C555, C556, C557, C558, C559, C560, C561, C562, C563, C564, C565, C566, C567, C568, C569, C570, C571, C572, C573, C574, C575, C576, C577, C578, C579, C580, C581, C582, C583, C584, C585, C586, C587, C588, C589, C590, C591, C592, C593, C594, C595, C596, C597, C598, C599, C600, C601, C602, C603, C604, C605, C606, C607, C608, C609, C610, C611, C612, C613, C614, C615, C616, C617, C618, C619, C620, C621, C622, C623, C624, C625, C626, C627, C628, C629, C630, C631, C632, C633, C634, C635, C636, C637, C638, C639, C640, C641, C642, C643, C644, C645, C646, C647, C648, C649, C650, C651, C652, C653, C654, C655, C656, C657, C658, C659, C660, C661, C662, C663, C664, C665, C666, C667, C668, C669, C670, C671, C672, C673, C674, C675, C676, C677, C678, C679, C680, C681, C682, C683, C684, C685, C686, C687, C688, C689, C690, C691, C692, C693, C694, C695, C696, C697, C698, C699, C700, C701, C702, C703, C704, C705, C706, C707, C708, C709, C710, C711, C712, C713, C714, C715, C716, C717, C718, C719, C720, C721, C722, C723, C724, C725, C726, C727, C728, C729, C730, C731, C732, C733, C734, C735, C736, C737, C738, C739, C740, C741, C742, C743, C744, C745, C746, C747, C748, C749, C750, C751, C752, C753, C754, C755, C756, C757, C758, C759, C760, C761, C762, C763, C764, C765, C766, C767, C768, C769, C770, C771, C772, C773, C774, C775, C776, C777, C778, C779, C780, C781, C782, C783, C784, C785, C786, C787, C788, C789, C790, C791, C792, C793, C794, C795, C796, C797, C798, C799, C800,

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
58	AA	1	Total 87	C 23	H 45	N 5	O 14	0	0
58	BA	1	Total 87	C 23	H 45	N 5	O 14	0	0
58	BA	1	Total 87	C 23	H 45	N 5	O 14	0	0
58	BA	1	Total 87	C 23	H 45	N 5	O 14	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
58	BA	1	Total C H N O 87 23 45 5 14	0	0
58	BA	1	Total C N O 42 23 5 14	0	0
58	CA	1	Total C N O 42 23 5 14	0	0
58	CA	1	Total C H N O 87 23 45 5 14	0	0
58	CA	1	Total C H N O 87 23 45 5 14	0	0
58	CA	1	Total C N O 42 23 5 14	0	0
58	CA	1	Total C H N O 87 23 45 5 14	0	0
58	DA	1	Total C H N O 87 23 45 5 14	0	0
58	DA	1	Total C H N O 87 23 45 5 14	0	0

- Molecule 59 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
59	B4	1	Total Zn 1 1	0	0
59	C4	1	Total Zn 1 1	0	0

- Molecule 60 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
60	AA	192	Total O 192 192	0	0
60	AE	3	Total O 3 3	0	0
60	AL	1	Total O 1 1	0	0
60	AN	2	Total O 2 2	0	0
60	AT	5	Total O 5 5	0	0
60	BA	626	Total O 626 626	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
60	BB	14	Total 14	O 14	0	0
60	BC	7	Total 7	O 7	0	0
60	BD	2	Total 2	O 2	0	0
60	BE	1	Total 1	O 1	0	0
60	BF	1	Total 1	O 1	0	0
60	BL	8	Total 8	O 8	0	0
60	BN	2	Total 2	O 2	0	0
60	BQ	1	Total 1	O 1	0	0
60	BS	1	Total 1	O 1	0	0
60	BT	3	Total 3	O 3	0	0
60	B3	1	Total 1	O 1	0	0
60	B4	1	Total 1	O 1	0	0
60	CA	608	Total 608	O 608	0	0
60	CB	14	Total 14	O 14	0	0
60	CC	10	Total 10	O 10	0	0
60	CD	5	Total 5	O 5	0	0
60	CE	3	Total 3	O 3	0	0
60	CJ	2	Total 2	O 2	0	0
60	CL	7	Total 7	O 7	0	0
60	CN	2	Total 2	O 2	0	0
60	CT	3	Total 3	O 3	0	0

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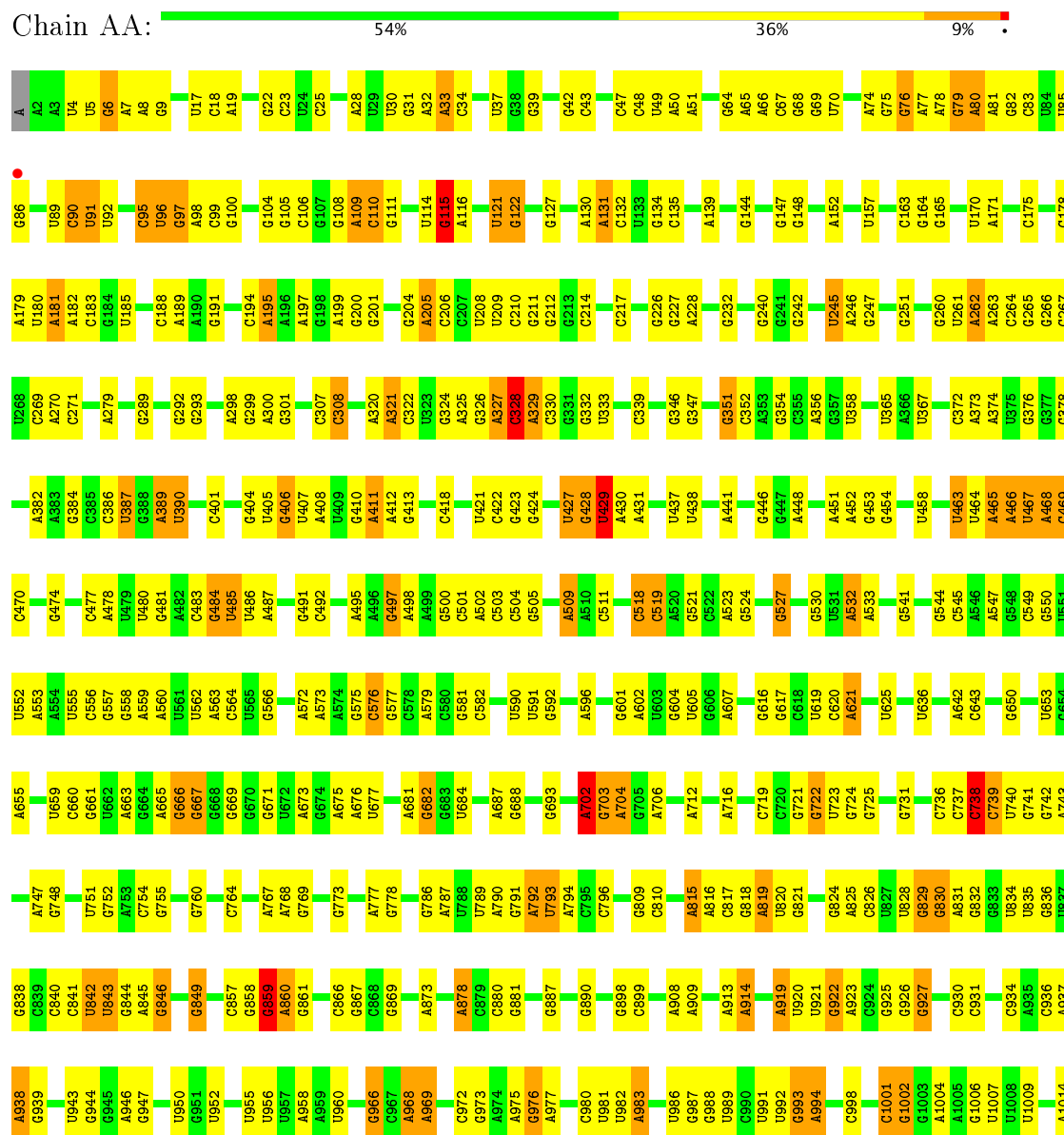
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
60	CU	1	Total 1	O 1	0	0
60	C3	1	Total 1	O 1	0	0
60	C4	2	Total 2	O 2	0	0
60	DA	184	Total 184	O 184	0	0
60	DD	2	Total 2	O 2	0	0
60	DK	2	Total 2	O 2	0	0
60	DL	2	Total 2	O 2	0	0
60	DN	4	Total 4	O 4	0	0
60	DT	3	Total 3	O 3	0	0
60	DU	1	Total 1	O 1	0	0

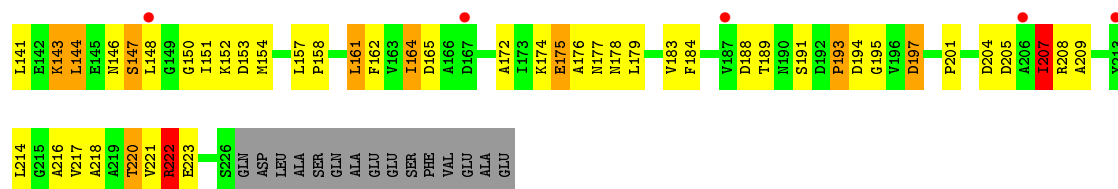
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: 16S ribosomal RNA

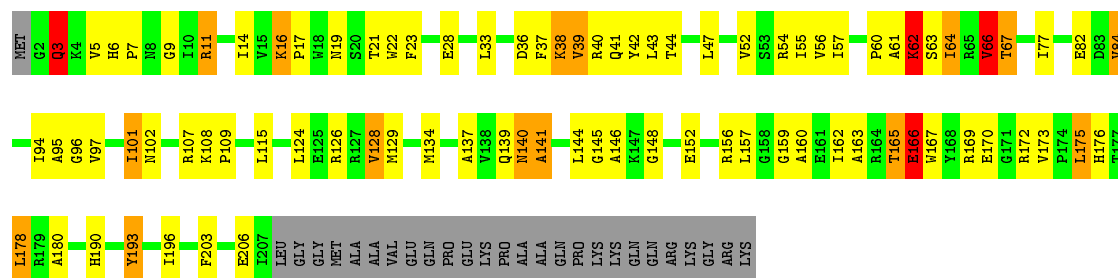






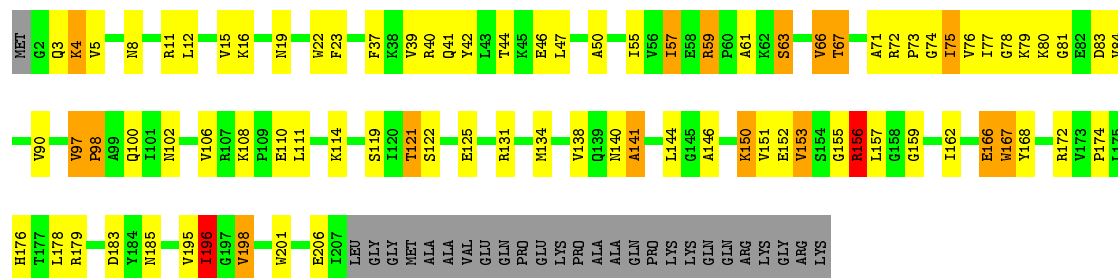
• Molecule 3: 30S ribosomal protein S3

Chain AC: 52% 29% 6% 12%



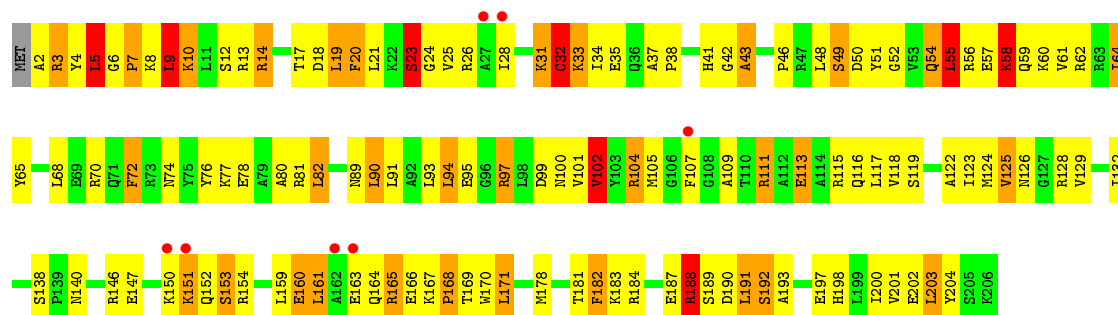
• Molecule 3: 30S ribosomal protein S3

Chain DC: 52% 29% 7% 12%



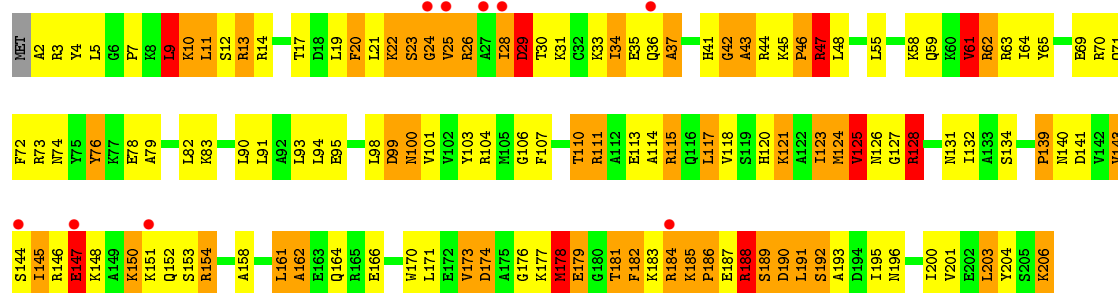
• Molecule 4: 30S ribosomal protein S4

Chain AD: 3% 37% 43% 16%

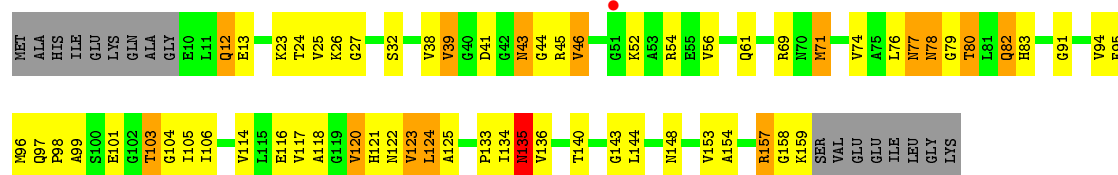


• Molecule 4: 30S ribosomal protein S4

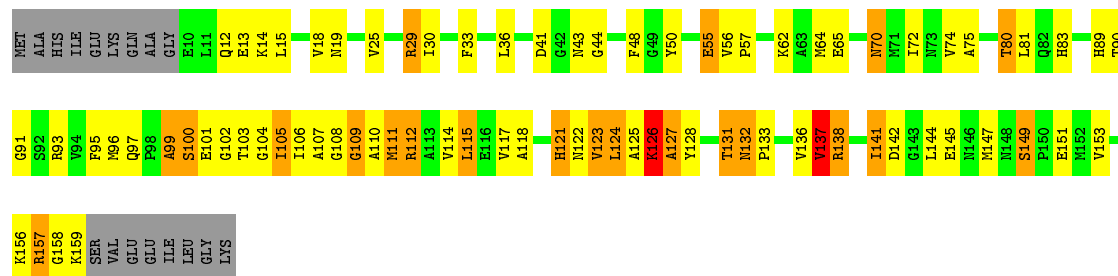
Chain DD: 4% 34% 38% 23%



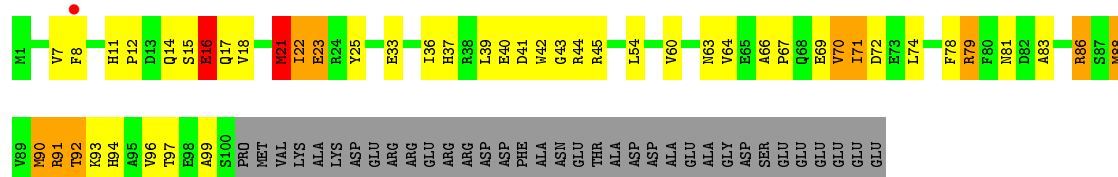
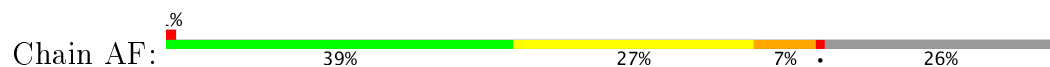
• Molecule 5: 30S ribosomal protein S5



• Molecule 5: 30S ribosomal protein S5



• Molecule 6: 30S ribosomal protein S6

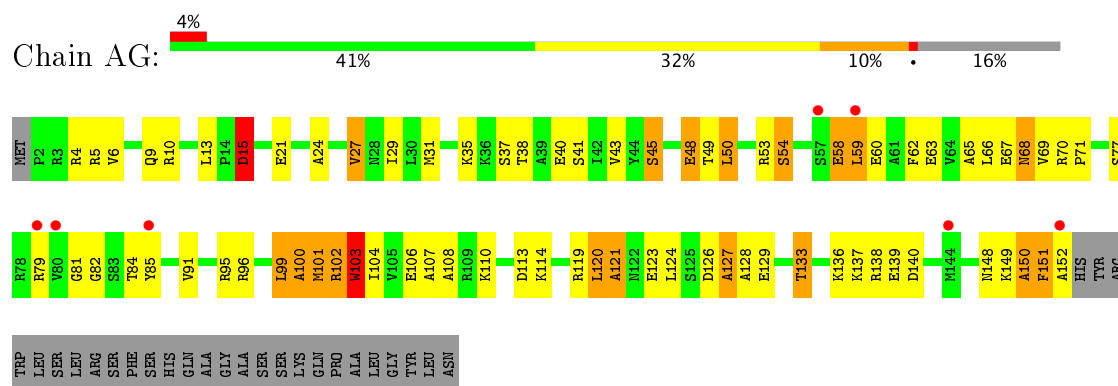


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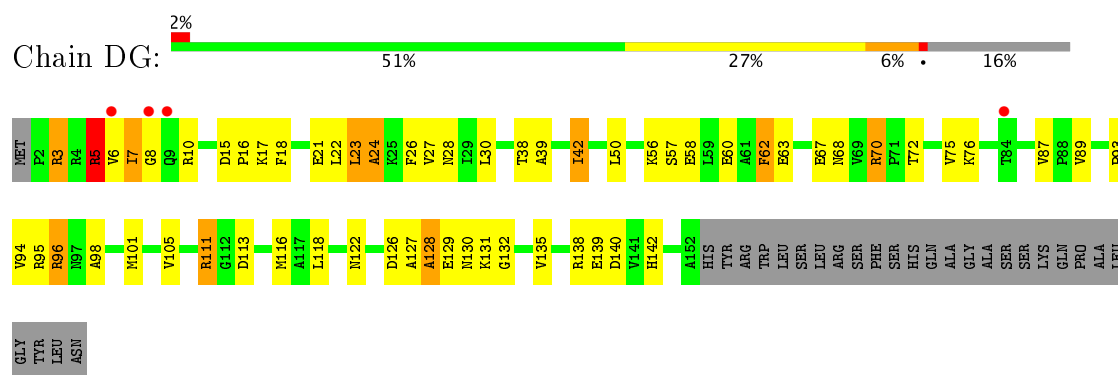




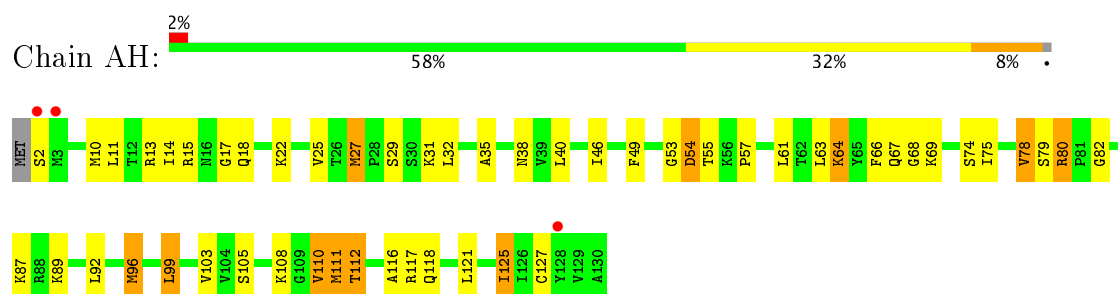
• Molecule 7: 30S ribosomal protein S7



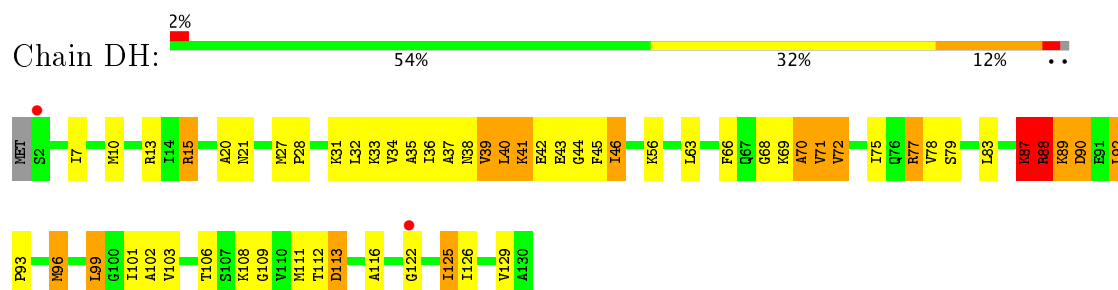
• Molecule 7: 30S ribosomal protein S7



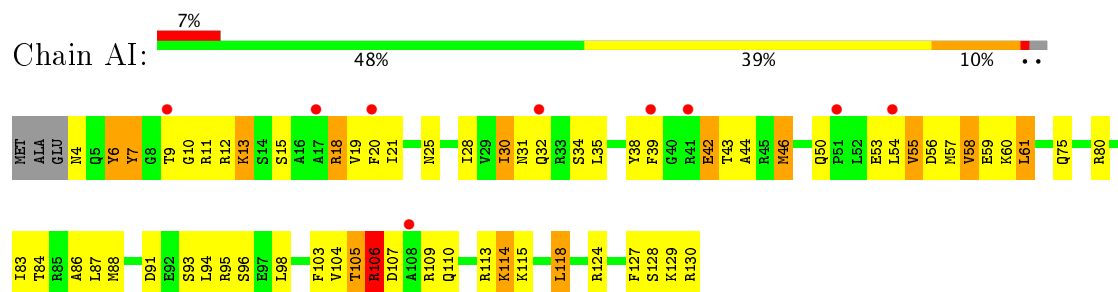
• Molecule 8: 30S ribosomal protein S8



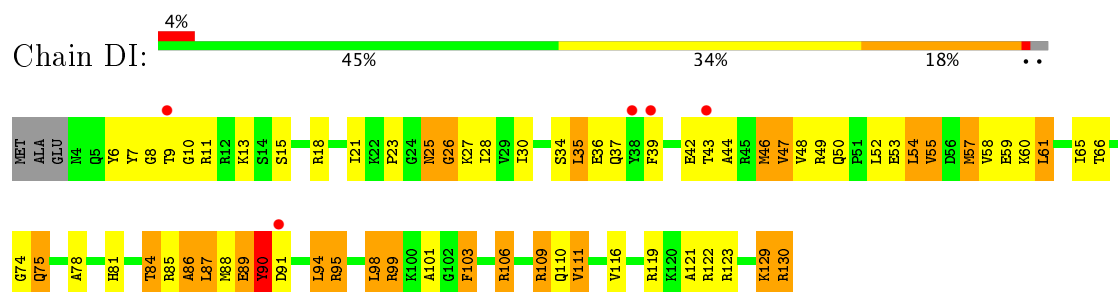
• Molecule 8: 30S ribosomal protein S8



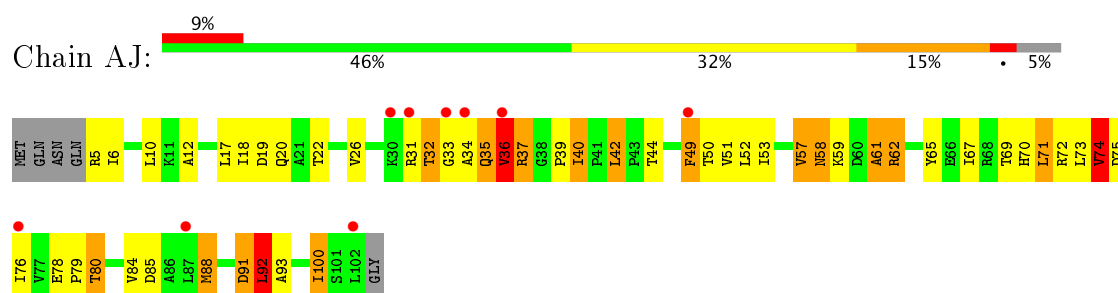
- Molecule 9: 30S ribosomal protein S9



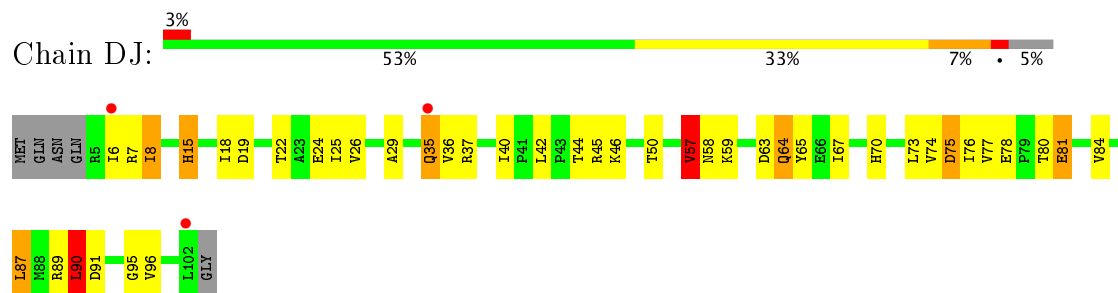
- Molecule 9: 30S ribosomal protein S9



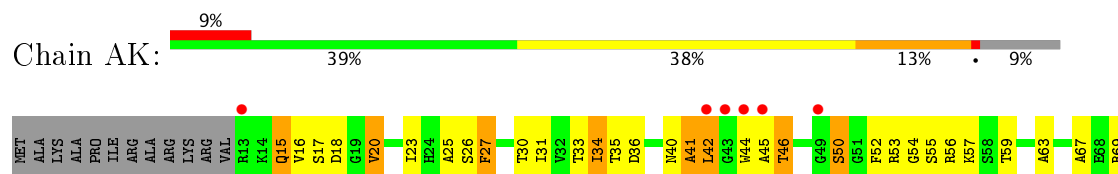
- Molecule 10: 30S ribosomal protein S10

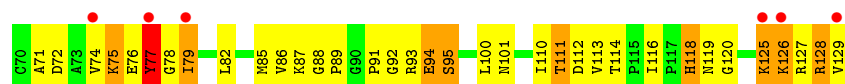


- Molecule 10: 30S ribosomal protein S10



- Molecule 11: 30S ribosomal protein S11

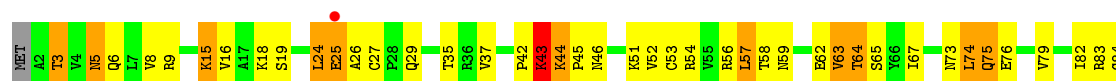




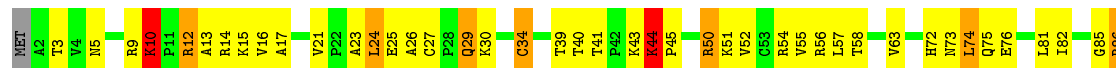
- Molecule 11: 30S ribosomal protein S11



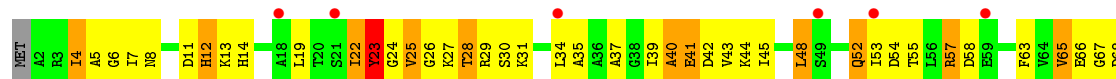
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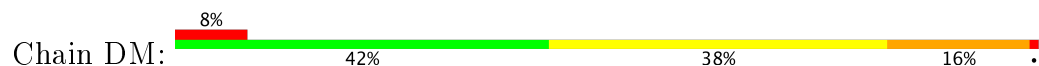
- Molecule 12: 30S ribosomal protein S12

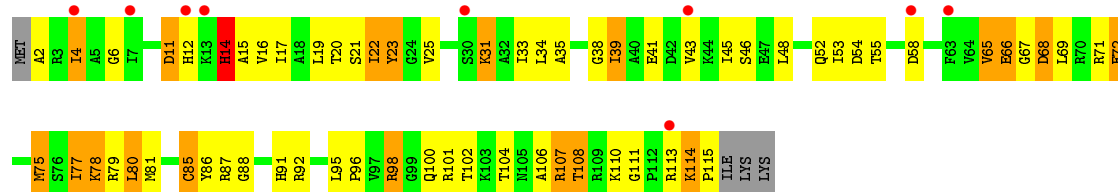


- Molecule 13: 30S ribosomal protein S13

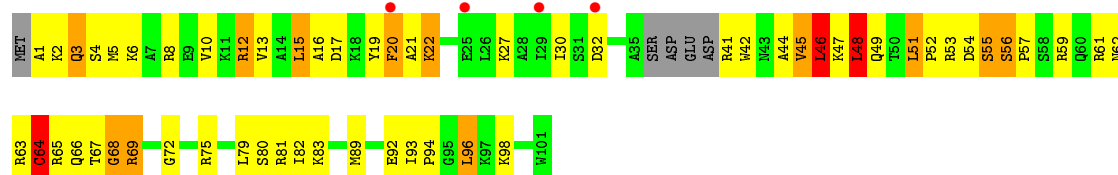


- Molecule 13: 30S ribosomal protein S13

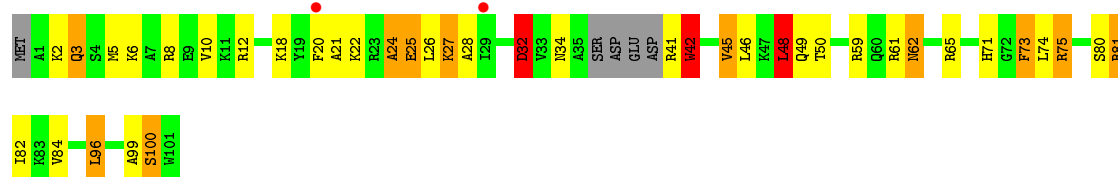




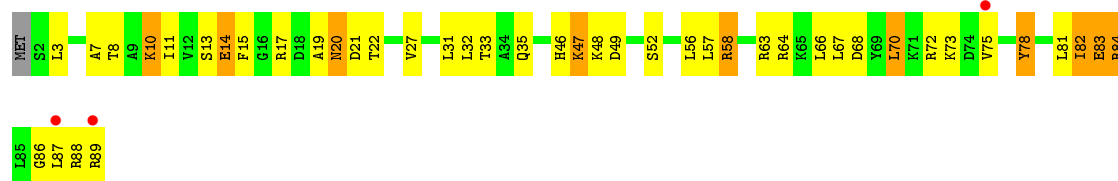
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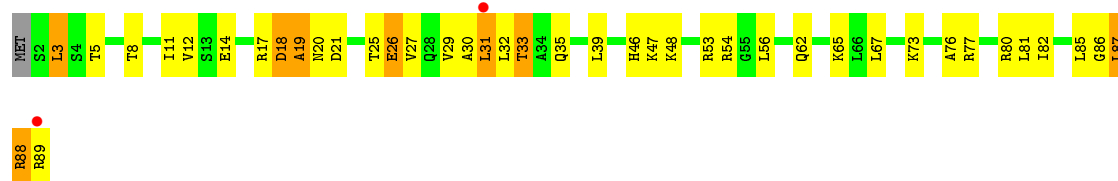
• Molecule 14: 30S ribosomal protein S14



• Molecule 15: 30S ribosomal protein S15

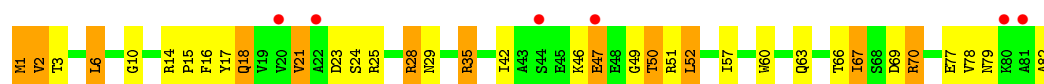


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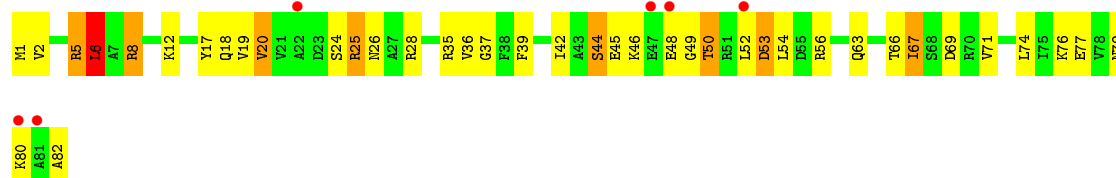


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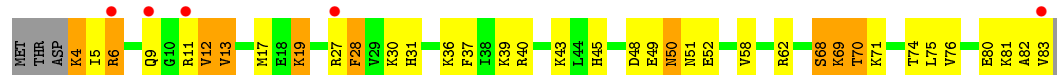




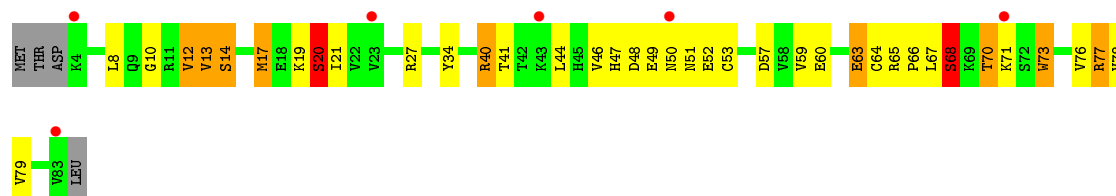
- Molecule 16: 30S ribosomal protein S16



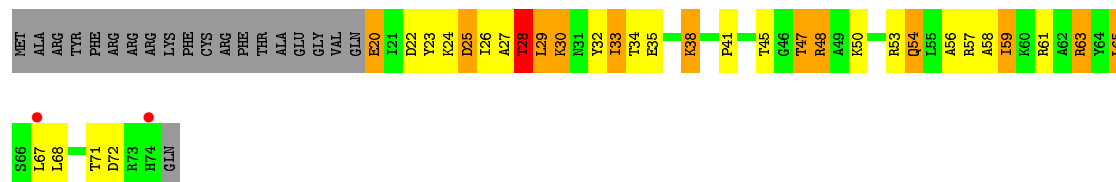
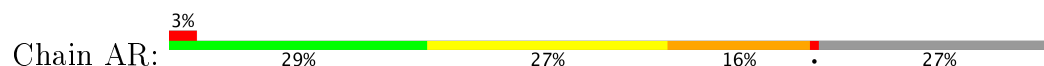
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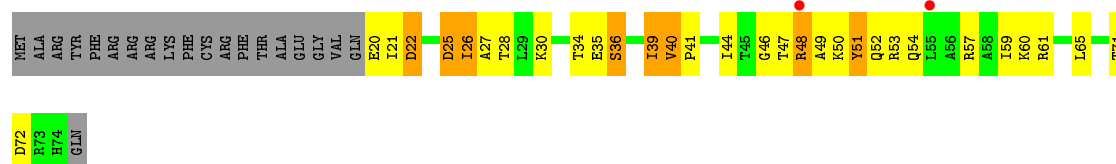
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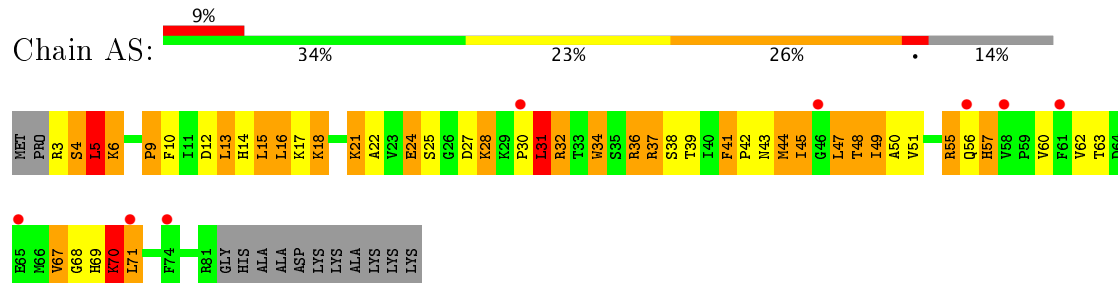
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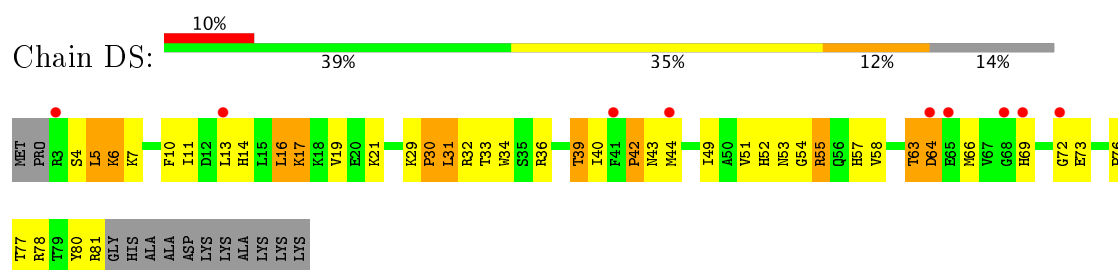
- Molecule 18: 30S ribosomal protein S18



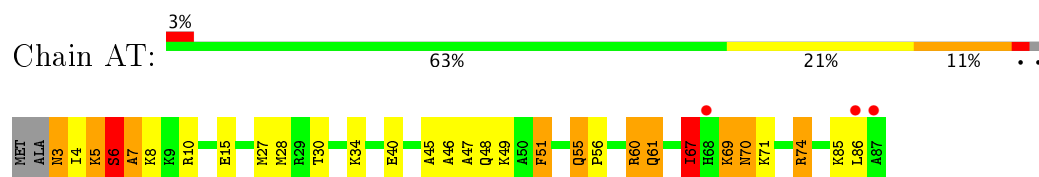
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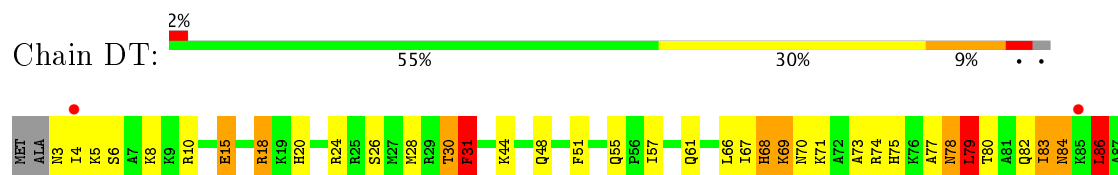
- Molecule 19: 30S ribosomal protein S19



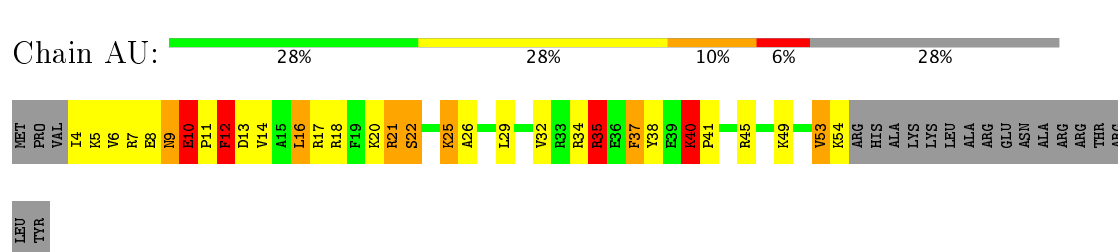
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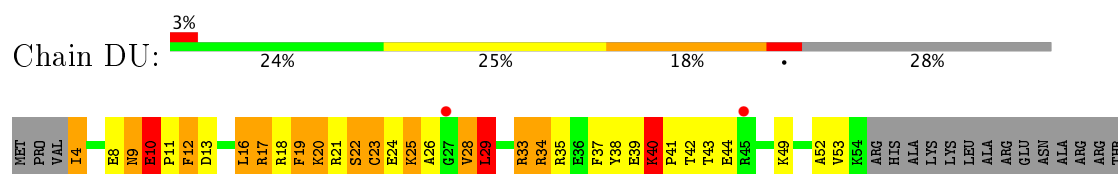
- Molecule 20: 30S ribosomal protein S20



- Molecule 21: 30S ribosomal protein S21

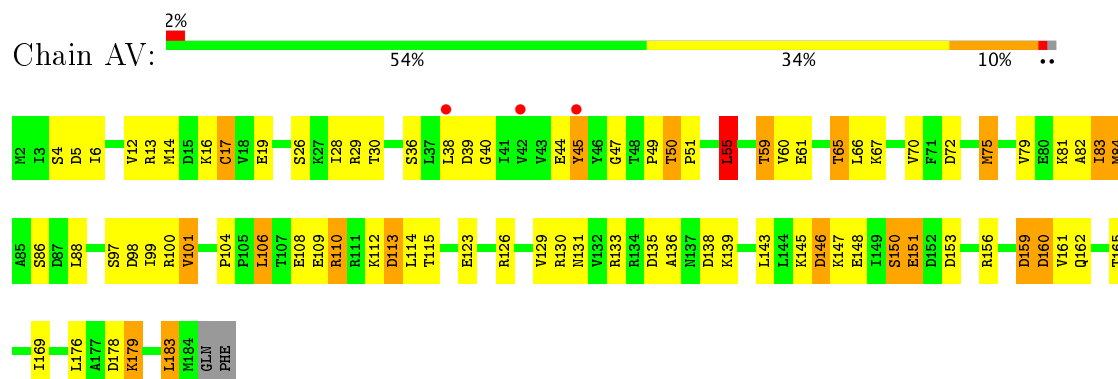


- Molecule 21: 30S ribosomal protein S21

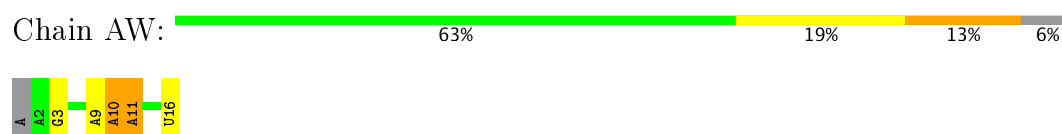


ARG
LEU
TYR

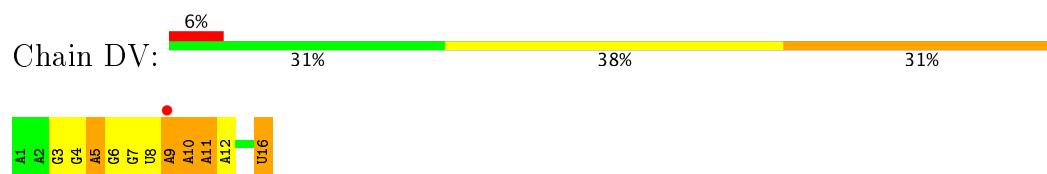
- Molecule 22: Ribosome-recycling factor



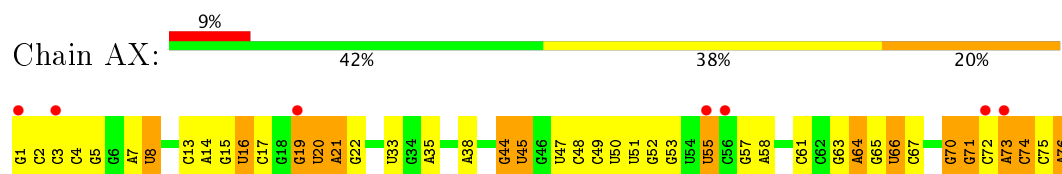
- Molecule 23: Messenger RNA



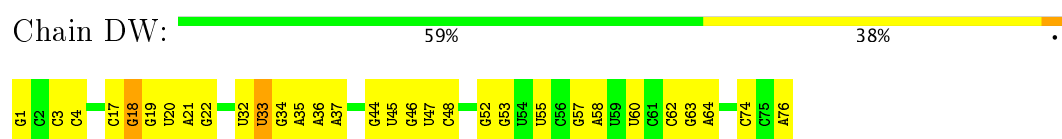
- Molecule 23: Messenger RNA



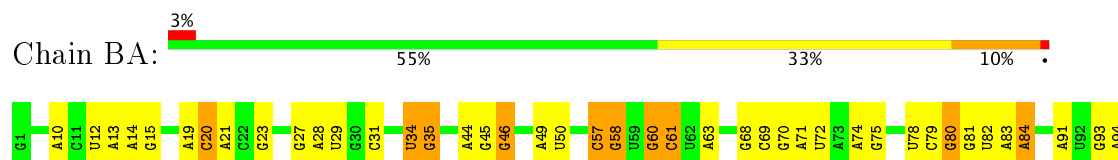
- Molecule 24: Phenylalanine specific transfer RNA, tRNA-Phe



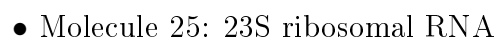
- Molecule 24: Phenylalanine specific transfer RNA, tRNA-Phe



- Molecule 25: 23S ribosomal RNA

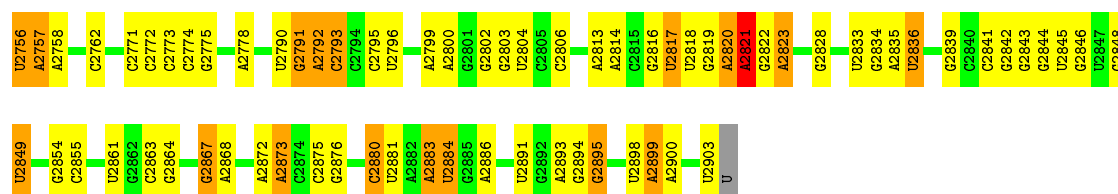


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	U1477	U1296	U1380	U1176	G1176	U1109	A1024	G923	A731				G450	A374	
	U1478	U1297	U1381	U1177	G1177	U1110	A1025	G924	A730				G451	A375	
	U1479	U1298	U1382	U1178	G1178	U1111	A1026	G925	A729				G452	A376	
	U1480	U1299	U1383	U1179	G1179	U1112	A1027	G926	A728				G453	A377	
	U1481	U1300	U1384	U1180	G1180	U1113	A1028	G927	A727				G454	A378	
	U1482	U1301	U1385	U1181	G1181	U1114	A1029	G928	A726				G455	A379	
	U1483	U1302	U1386	U1182	G1182	U1115	A1030	G929	A725				G456	A380	
	U1484	U1303	U1387	U1183	G1183	U1116	A1031	G930	A724				G457	A381	
	U1485	U1304	U1388	U1184	G1184	U1117	A1032	G931	A723				G458	A382	
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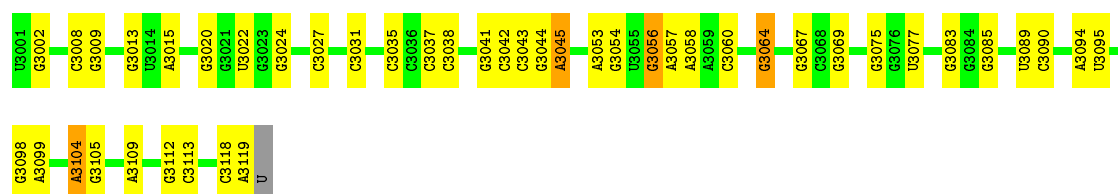


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	G2661	C2575	U2408	U2322	U2322		G2158	U2092	U2011	C1914	U1827	A1735	A1610	
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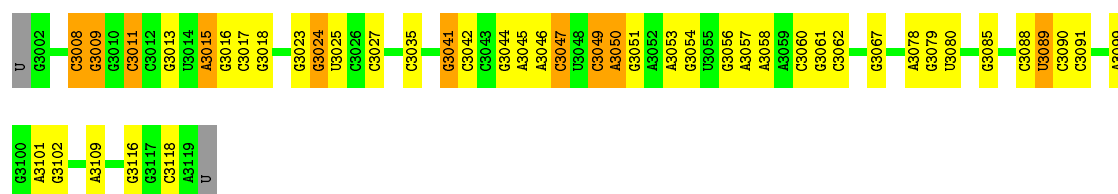
• Molecule 26: 5S ribosomal RNA

Chain BB: 63% 33% ..



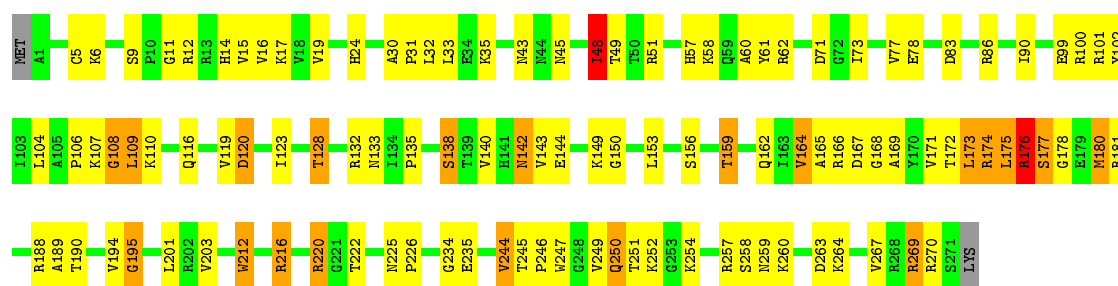
• Molecule 26: 5S ribosomal RNA

Chain CB: 61% 29% 8% .



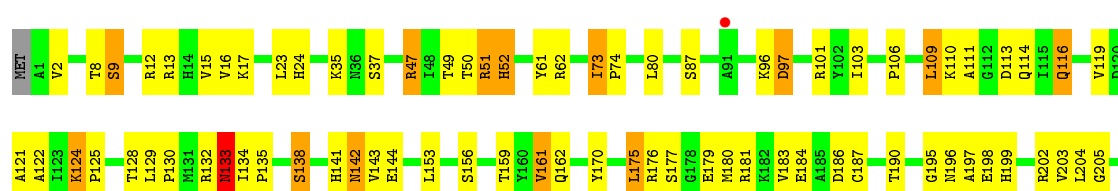
• Molecule 27: 50S ribosomal protein L2

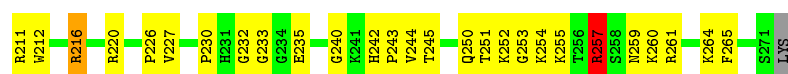
Chain BC: 59% 33% 7% ..



• Molecule 27: 50S ribosomal protein L2

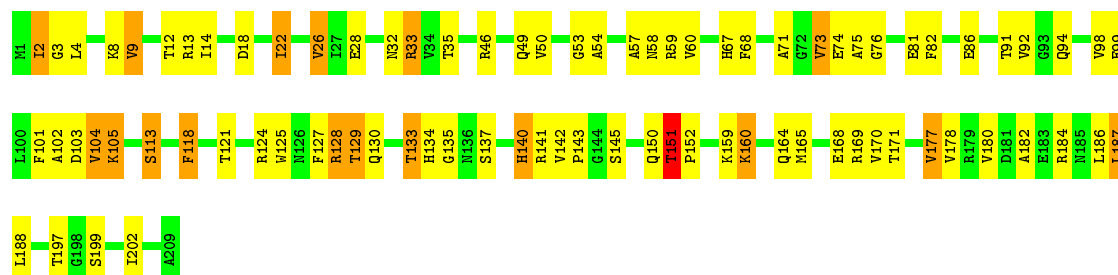
Chain CC: 61% 32% 5% ..





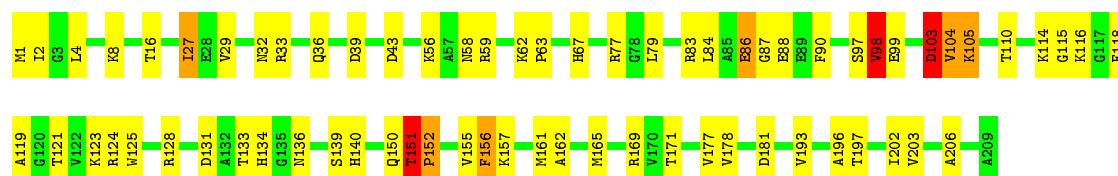
- Molecule 28: 50S ribosomal protein L3

Chain BD: 60% 32% 8%



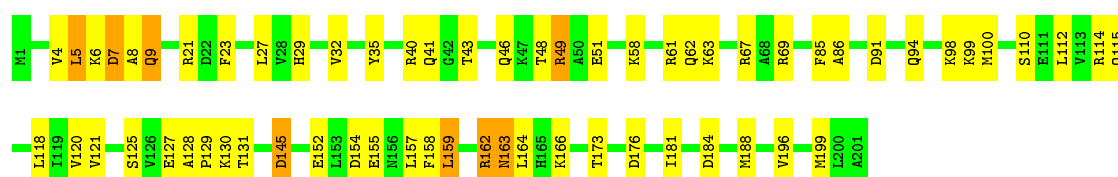
- Molecule 28: 50S ribosomal protein L3

Chain CD: 67% 29% 4%



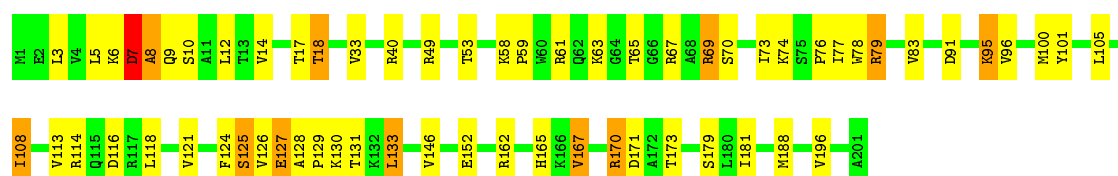
- Molecule 29: 50S ribosomal protein L4

Chain BE: 69% 27% 4%



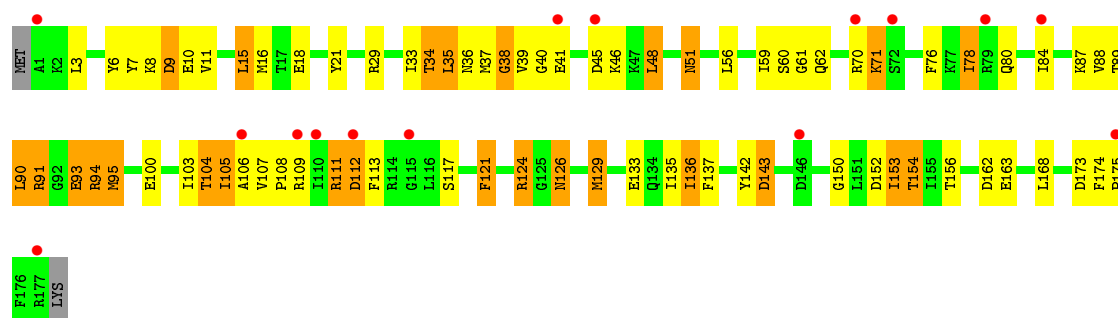
- Molecule 29: 50S ribosomal protein L4

Chain CE: 69% 25% 6%

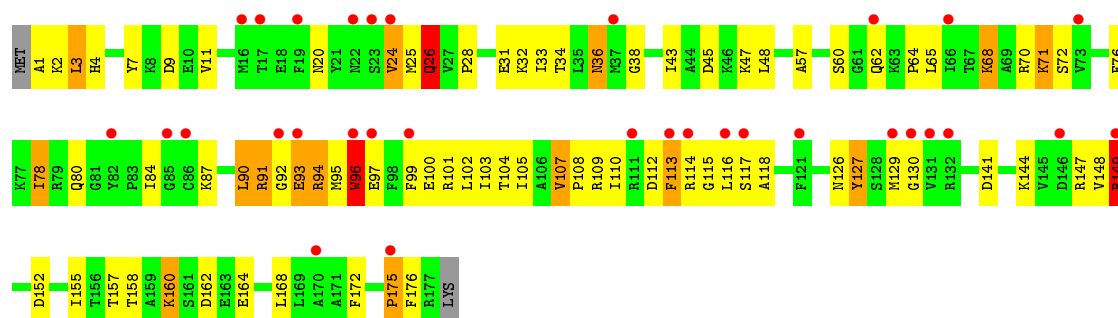


- Molecule 30: 50S ribosomal protein L5

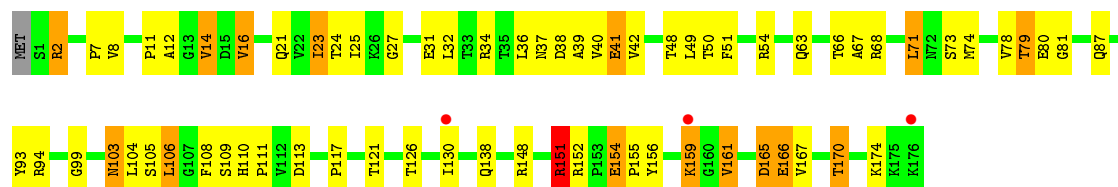
Chain BF: 8% 56% 28% 8%



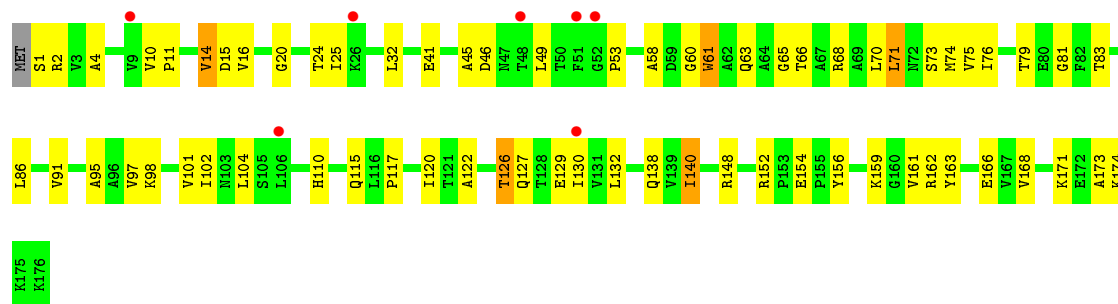
- Molecule 30: 50S ribosomal protein L5



- Molecule 31: 50S ribosomal protein L6

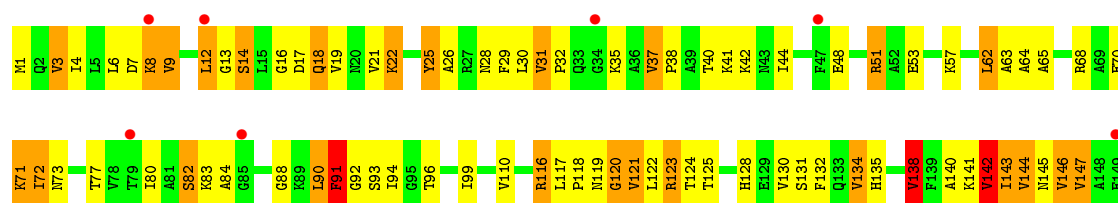


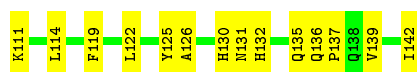
- Molecule 31: 50S ribosomal protein L6



- Molecule 32: 50S ribosomal protein L9

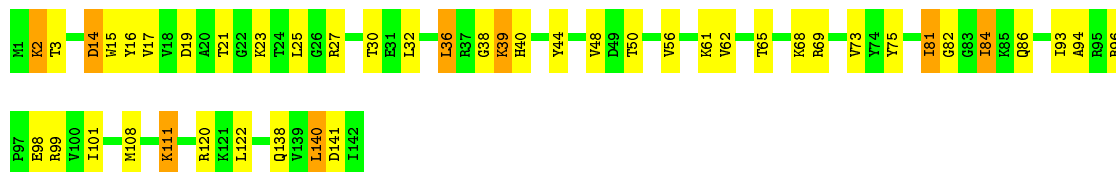






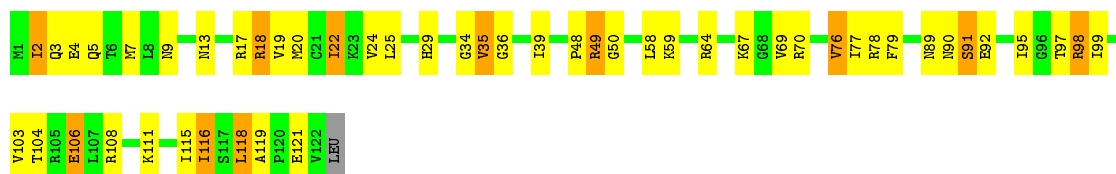
• Molecule 34: 50S ribosomal protein L13

Chain CJ: 68% 26% 6%



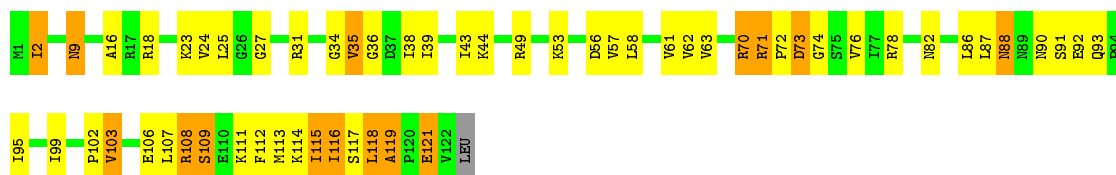
• Molecule 35: 50S ribosomal protein L14

Chain BK: 59% 32% 9%



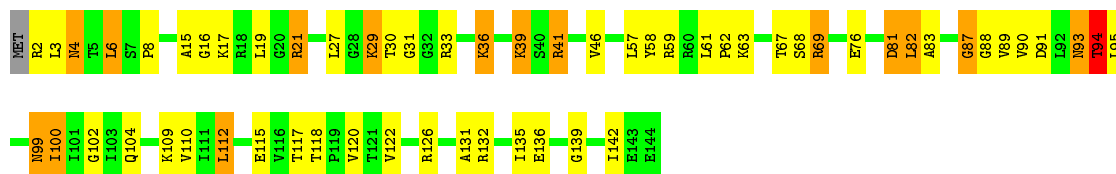
• Molecule 35: 50S ribosomal protein L14

Chain CK: 53% 34% 12%



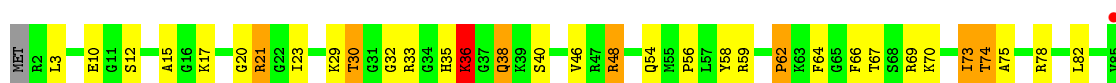
• Molecule 36: 50S ribosomal protein L15

Chain BL: 58% 30% 10%



• Molecule 36: 50S ribosomal protein L15

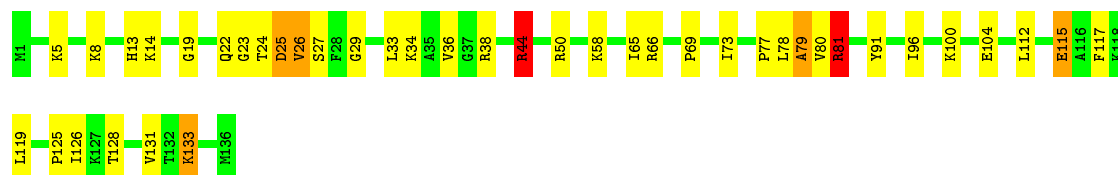
Chain CL: 56% 33% 10%





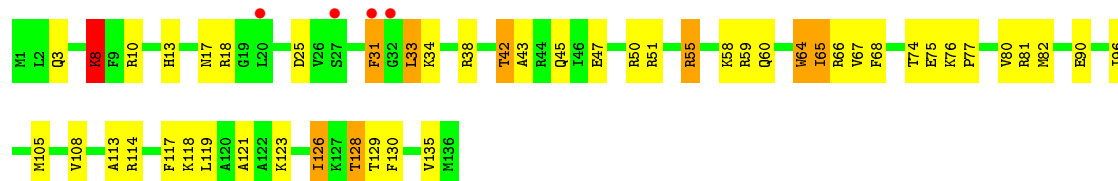
- Molecule 37: 50S ribosomal protein L16

Chain BM: 70% 25% . .



- Molecule 37: 50S ribosomal protein L16

Chain CM: 3% 64% 29% 6% .



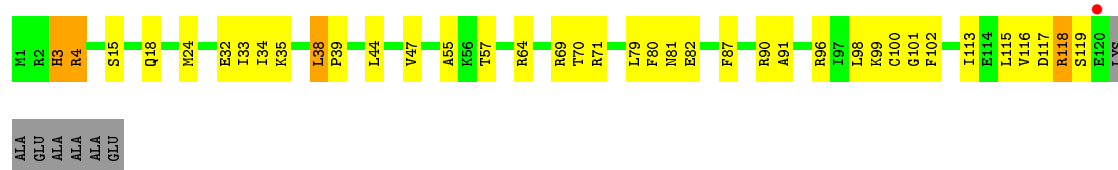
- Molecule 38: 50S ribosomal protein L17

Chain BN: 56% 30% 8% 6% .



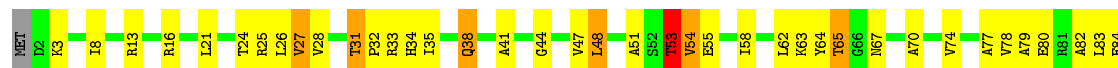
- Molecule 38: 50S ribosomal protein L17

Chain CN: .% 65% 27% 6% .



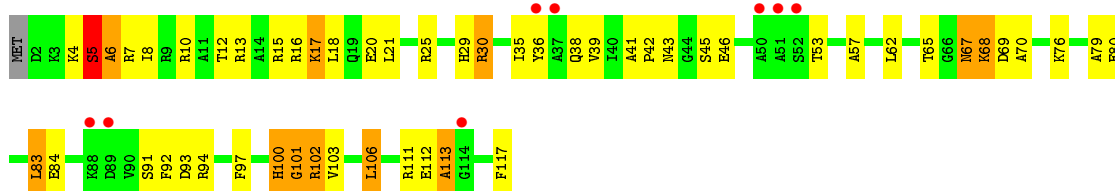
- Molecule 39: 50S ribosomal protein L18

Chain BO: 57% 35% 6% . .

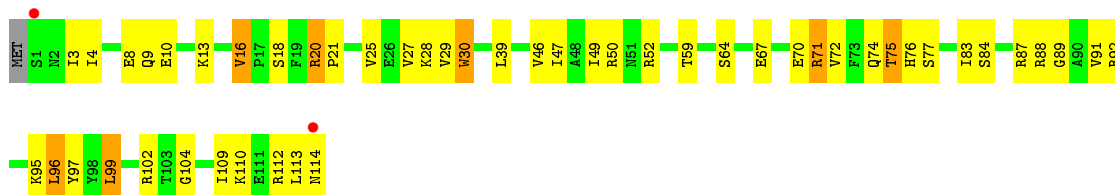




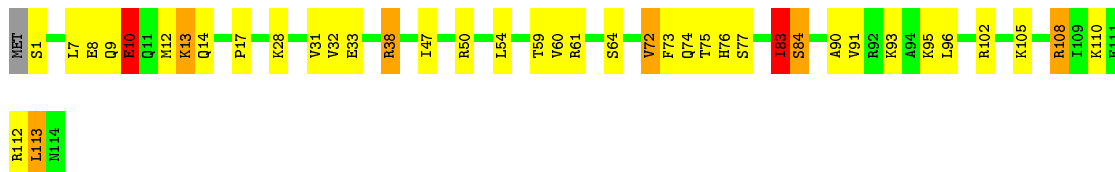
- Molecule 39: 50S ribosomal protein L18



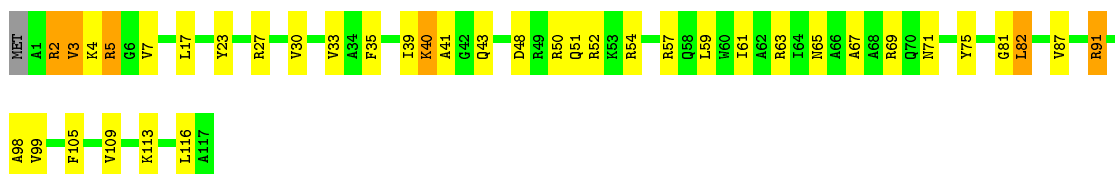
- Molecule 40: 50S ribosomal protein L19



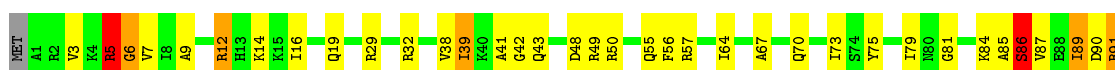
- Molecule 40: 50S ribosomal protein L19



- Molecule 41: 50S ribosomal protein L20



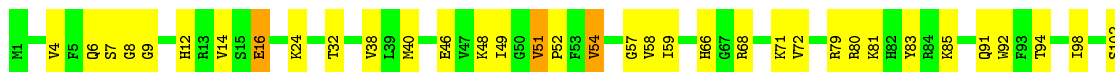
- Molecule 41: 50S ribosomal protein L20





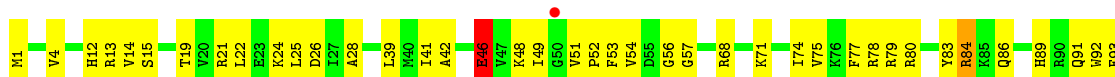
- Molecule 42: 50S ribosomal protein L21

Chain BR: 66% 31%



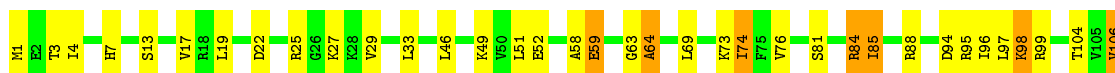
- Molecule 42: 50S ribosomal protein L21

Chain CR: 60% 38%



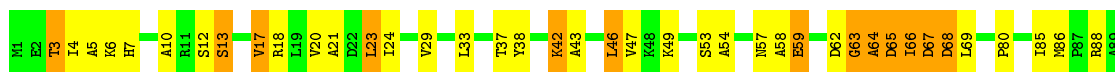
- Molecule 43: 50S ribosomal protein L22

Chain BS: 65% 29% 6%



- Molecule 43: 50S ribosomal protein L22

Chain CS: 55% 33% 13%



- Molecule 44: 50S ribosomal protein L23

Chain BT: 56% 25% 12% 7%



GLU

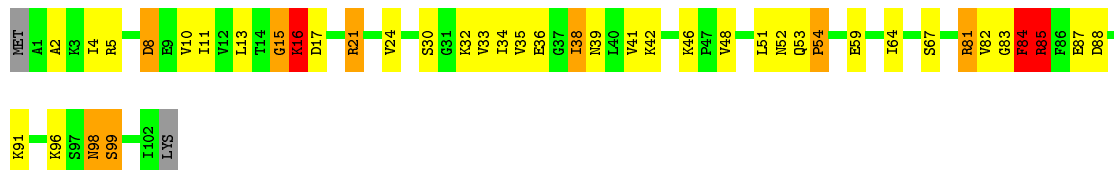
- Molecule 44: 50S ribosomal protein L23

Chain CT: 



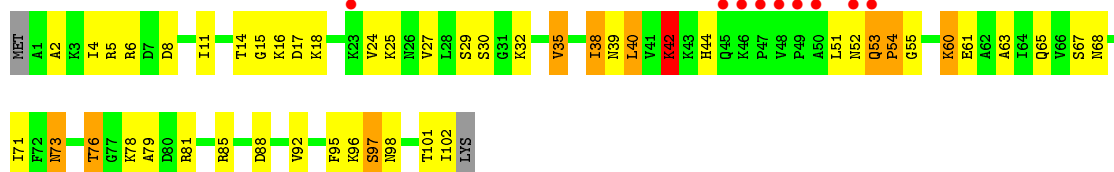
- Molecule 45: 50S ribosomal protein L24

Chain BU: 




- Molecule 45: 50S ribosomal protein L24

Chain CU: 




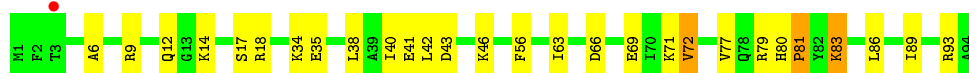
- Molecule 46: 50S ribosomal protein L25

Chain BV: 



- Molecule 46: 50S ribosomal protein L25

Chain CV: 

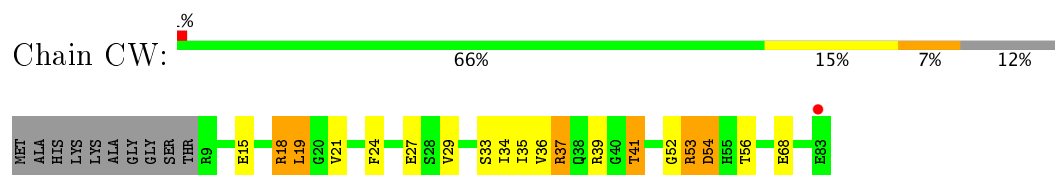


- Molecule 47: 50S ribosomal protein L27

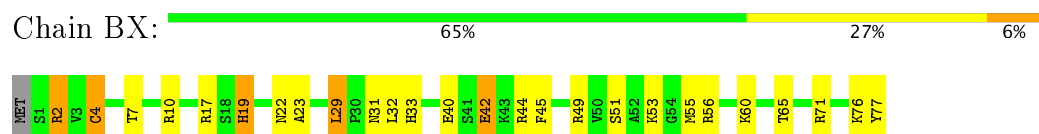
Chain BW: 



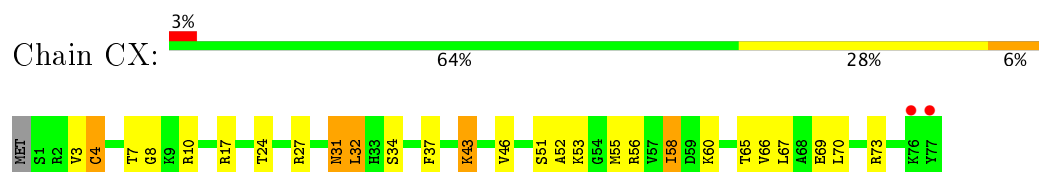
- Molecule 47: 50S ribosomal protein L27



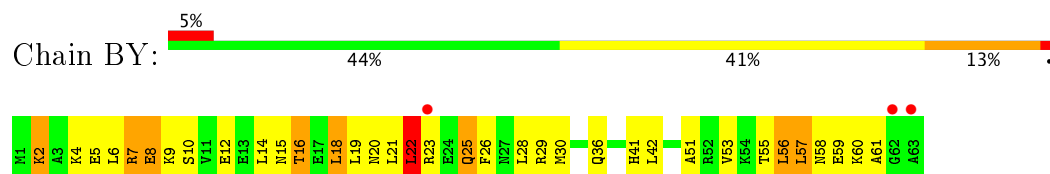
- Molecule 48: 50S ribosomal protein L28



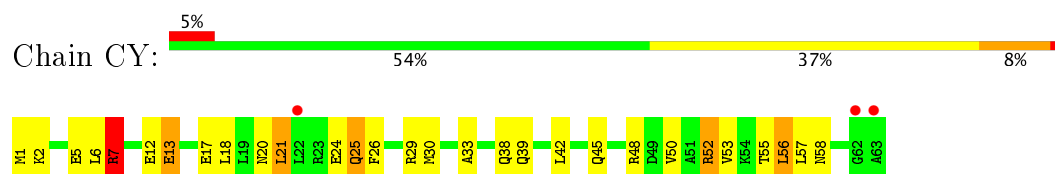
- Molecule 48: 50S ribosomal protein L28



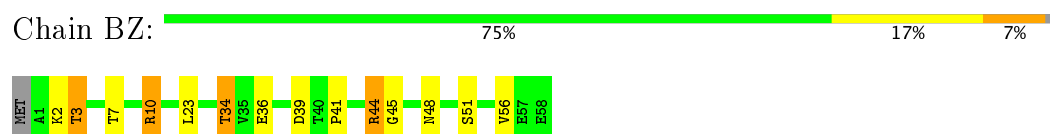
- Molecule 49: 50S ribosomal protein L29



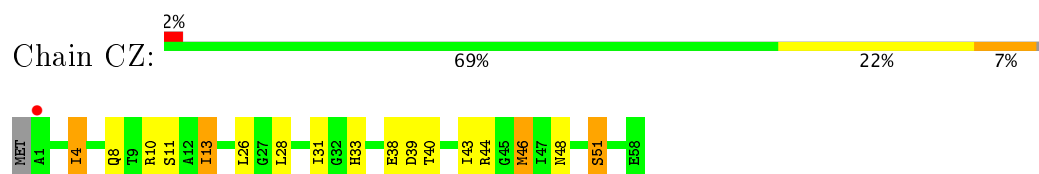
- Molecule 49: 50S ribosomal protein L29



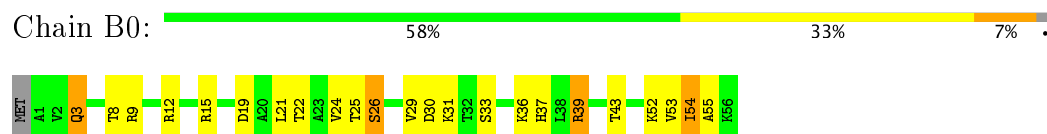
- Molecule 50: 50S ribosomal protein L30



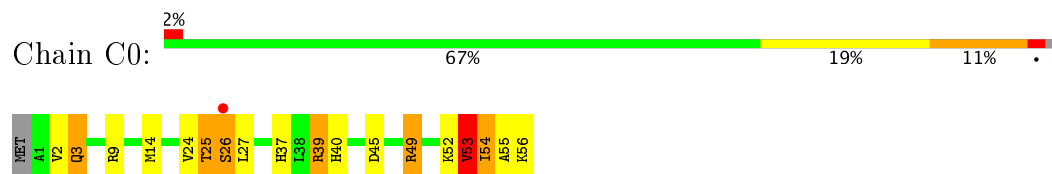
- Molecule 50: 50S ribosomal protein L30



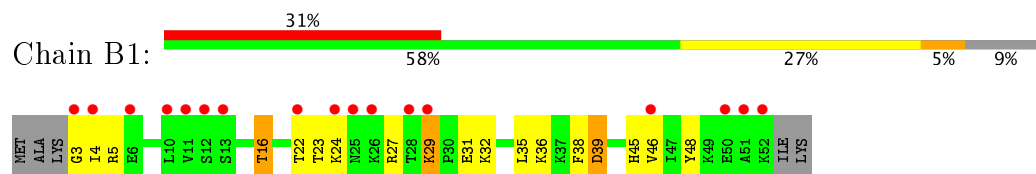
- Molecule 51: 50S ribosomal protein L32



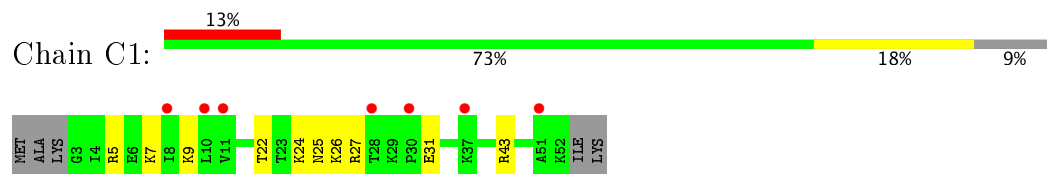
- Molecule 51: 50S ribosomal protein L32



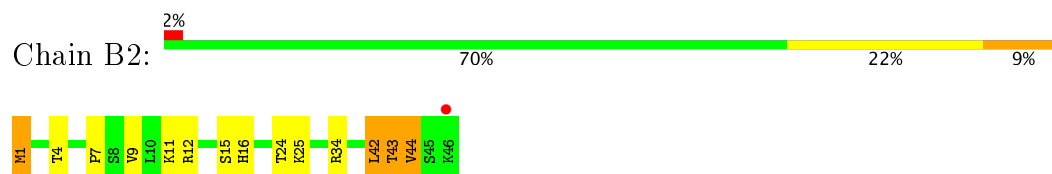
- Molecule 52: 50S ribosomal protein L33



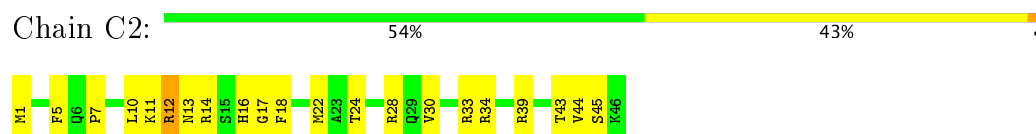
- Molecule 52: 50S ribosomal protein L33



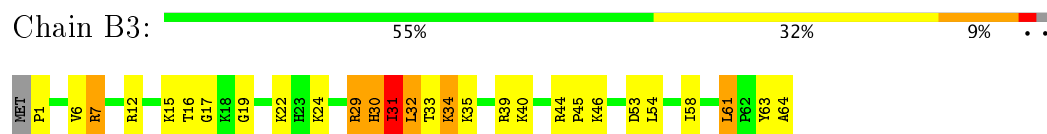
- Molecule 53: 50S ribosomal protein L34



- Molecule 53: 50S ribosomal protein L34



- Molecule 54: 50S ribosomal protein L35



- Molecule 54: 50S ribosomal protein L35





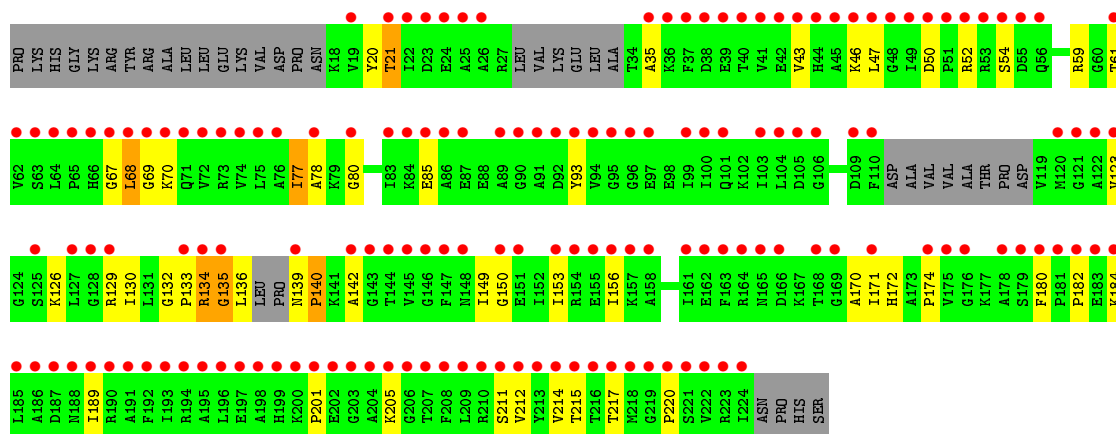
- Molecule 55: 50S ribosomal protein L36



- Molecule 55: 50S ribosomal protein L36



- Molecule 56: 50S ribosomal protein L1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	212.10Å 435.24Å 614.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.00 – 3.00 69.34 – 2.93	Depositor EDS
% Data completeness (in resolution range)	81.1 (70.00-3.00) 76.0 (69.34-2.93)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.91Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.174 , 0.253 0.178 , 0.257	Depositor DCC
R_{free} test set	4548 reflections (0.50%)	DCC
Wilson B-factor (Å ²)	72.9	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 95.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	294484	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, PAR

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	AA	0.48	0/36944	0.99	30/57632 (0.1%)
1	DA	0.49	0/36966	1.00	37/57666 (0.1%)
2	AB	0.38	0/1735	0.64	0/2338
2	DB	0.41	0/1735	0.67	1/2338 (0.0%)
3	AC	0.45	0/1651	0.70	0/2225
3	DC	0.42	0/1651	0.66	0/2225
4	AD	0.46	0/1665	0.75	3/2227 (0.1%)
4	DD	0.51	0/1665	0.81	1/2227 (0.0%)
5	AE	0.40	0/1118	0.62	0/1504
5	DE	0.47	0/1118	0.73	0/1504
6	AF	0.42	0/835	0.69	0/1128
6	DF	0.43	0/835	0.71	0/1128
7	AG	2.14	8/1195 (0.7%)	0.81	4/1602 (0.2%)
7	DG	0.38	0/1195	0.57	0/1602
8	AH	0.41	0/989	0.66	0/1326
8	DH	0.45	0/989	0.77	1/1326 (0.1%)
9	AI	0.38	0/1034	0.65	0/1375
9	DI	0.42	0/1034	0.66	0/1375
10	AJ	0.42	0/796	0.69	1/1077 (0.1%)
10	DJ	0.38	0/796	0.64	0/1077
11	AK	0.43	0/893	0.69	0/1205
11	DK	0.43	0/893	0.67	0/1205
12	AL	0.52	0/969	0.77	0/1300
12	DL	0.52	0/969	0.81	1/1300 (0.1%)
13	AM	0.39	0/892	0.74	0/1193
13	DM	0.36	0/892	0.66	0/1193
14	AN	0.44	0/785	0.73	0/1043
14	DN	0.39	0/785	0.67	0/1043
15	AO	0.40	0/724	0.67	0/966
15	DO	0.38	0/724	0.63	0/966
16	AP	0.40	0/659	0.62	0/884
16	DP	0.44	0/659	0.72	1/884 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AQ	0.40	0/657	0.72	0/881
17	DQ	0.45	0/657	0.69	0/881
18	AR	0.44	0/462	0.63	0/621
18	DR	0.44	0/462	0.71	1/621 (0.2%)
19	AS	0.37	0/652	0.71	0/877
19	DS	0.37	0/652	0.61	0/877
20	AT	0.41	0/671	0.64	0/888
20	DT	0.41	0/671	0.64	0/888
21	AU	0.50	0/430	0.91	1/570 (0.2%)
21	DU	0.58	0/430	0.89	0/570
22	AV	0.43	0/1430	0.63	0/1924
23	AW	0.58	0/363	1.12	0/564
23	DV	0.51	0/388	1.04	0/603
24	AX	0.49	1/1813 (0.1%)	0.95	0/2823
24	DW	0.52	1/1813 (0.1%)	0.95	0/2823
25	BA	0.67	8/69659 (0.0%)	1.20	303/108672 (0.3%)
25	CA	0.52	3/69659 (0.0%)	1.04	112/108672 (0.1%)
26	BB	0.54	0/2850	1.03	7/4444 (0.2%)
26	CB	0.40	0/2828	0.90	2/4410 (0.0%)
27	BC	0.47	0/2121	0.71	0/2852
27	CC	0.42	0/2121	0.68	0/2852
28	BD	0.53	0/1586	0.77	2/2134 (0.1%)
28	CD	0.44	0/1586	0.73	1/2134 (0.0%)
29	BE	0.49	0/1571	0.71	0/2113
29	CE	0.43	0/1571	0.65	0/2113
30	BF	0.40	0/1434	0.65	0/1926
30	CF	0.36	0/1434	0.61	0/1926
31	BG	0.43	0/1343	0.61	0/1816
31	CG	0.35	0/1343	0.56	0/1816
32	BH	0.43	0/1118	0.70	0/1511
32	CH	0.38	0/1118	0.60	1/1511 (0.1%)
33	BI	0.38	0/1046	0.60	0/1410
33	CI	0.33	0/1046	0.55	0/1410
34	BJ	0.53	0/1152	0.67	0/1551
34	CJ	0.43	0/1152	0.62	0/1551
35	BK	0.54	0/947	0.74	0/1268
35	CK	0.47	0/947	0.72	0/1268
36	BL	0.50	0/1054	0.80	0/1403
36	CL	0.45	0/1054	0.70	0/1403
37	BM	0.50	0/1093	0.75	1/1460 (0.1%)
37	CM	0.42	0/1093	0.59	0/1460
38	BN	0.49	0/973	0.69	0/1301
38	CN	0.50	0/973	0.66	1/1301 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
39	BO	0.45	0/902	0.65	0/1209
39	CO	0.37	0/902	0.65	0/1209
40	BP	0.49	0/929	0.66	0/1242
40	CP	0.45	0/929	0.65	1/1242 (0.1%)
41	BQ	0.53	0/960	0.74	0/1278
41	CQ	0.49	0/960	0.72	1/1278 (0.1%)
42	BR	0.52	0/829	0.72	0/1107
42	CR	0.49	0/829	0.71	0/1107
43	BS	0.53	0/864	0.70	0/1156
43	CS	0.47	0/864	0.75	0/1156
44	BT	0.48	0/744	0.67	0/994
44	CT	0.40	0/744	0.56	0/994
45	BU	0.51	0/787	0.72	0/1051
45	CU	0.43	0/787	0.64	0/1051
46	BV	0.45	0/766	0.64	0/1025
46	CV	0.34	0/766	0.58	0/1025
47	BW	0.54	0/587	0.73	0/776
47	CW	0.41	0/576	0.65	0/762
48	BX	0.47	0/635	0.71	0/848
48	CX	0.38	0/635	0.64	0/848
49	BY	0.42	0/510	0.69	0/677
49	CY	0.40	0/510	0.72	0/677
50	BZ	0.51	0/453	0.78	0/605
50	CZ	0.41	0/453	0.62	0/605
51	B0	0.54	0/450	0.79	1/599 (0.2%)
51	C0	0.48	0/450	0.73	0/599
52	B1	0.39	0/416	0.61	0/554
52	C1	0.38	0/416	0.62	0/554
53	B2	0.50	0/380	0.74	0/498
53	C2	0.45	0/380	0.66	0/498
54	B3	0.55	0/513	0.78	0/676
54	C3	0.46	0/513	0.71	0/676
55	B4	0.53	0/303	0.77	0/397
55	C4	0.42	0/303	0.60	0/397
56	B5	0.33	0/1145	0.54	0/1556
All	All	0.54	21/316403 (0.0%)	0.99	515/473109 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	AB	0	1
2	DB	0	1
4	AD	0	1
4	DD	0	1
5	DE	0	2
6	AF	0	1
7	AG	0	1
9	DI	0	1
10	AJ	0	2
11	AK	0	1
12	DL	0	1
14	AN	0	2
17	DQ	0	1
19	AS	0	1
20	AT	0	1
20	DT	0	2
21	AU	0	1
27	BC	0	3
28	BD	0	1
28	CD	0	2
30	BF	0	1
36	BL	0	1
36	CL	0	1
38	BN	0	1
38	CN	0	1
40	CP	0	1
42	CR	0	1
44	BT	0	1
48	CX	0	1
All	All	0	36

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	AG	103	TRP	CE3-CZ3	40.14	2.06	1.38
7	AG	100	ALA	CA-C	39.86	2.56	1.52
7	AG	103	TRP	CZ3-CH2	28.41	1.85	1.40
7	AG	103	TRP	CE2-CZ2	21.77	1.76	1.39
7	AG	103	TRP	CD2-CE2	18.20	1.63	1.41

The worst 5 of 515 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	BA	984	A	C2-N3-C4	-15.53	102.84	110.60
25	BA	1936	A	C2-N3-C4	-12.94	104.13	110.60
28	CD	151	THR	C-N-CD	-11.23	95.90	120.60
25	BA	2623	G	O5'-P-OP1	-10.98	95.82	105.70
25	BA	1936	A	N1-C2-N3	10.97	134.79	129.30

There are no chirality outliers.

5 of 36 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	AB	148	LEU	Peptide
4	AD	18	ASP	Peptide
6	AF	79	ARG	Peptide
7	AG	126	ASP	Peptide
10	AJ	17	LEU	Peptide

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	AA	32995	0	16607	432	1
1	DA	33015	0	16617	426	0
2	AB	1704	0	1732	37	0
2	DB	1704	0	1732	80	0
3	AC	1624	0	1696	45	0
3	DC	1624	0	1696	39	0
4	AD	1643	0	1707	92	0
4	DD	1643	0	1707	117	0
5	AE	1105	0	1148	32	0
5	DE	1105	0	1148	53	0
6	AF	817	0	808	20	0
6	DF	817	0	808	22	0
7	AG	1181	0	1238	78	0
7	DG	1181	0	1238	25	0
8	AH	979	0	1031	34	0
8	DH	979	0	1031	49	0
9	AI	1022	0	1070	36	0
9	DI	1022	0	1070	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	AJ	786	0	828	36	0
10	DJ	786	0	828	16	0
11	AK	877	0	887	46	0
11	DK	877	0	887	35	0
12	AL	955	0	1016	47	0
12	DL	955	0	1016	42	0
13	AM	883	0	941	62	0
13	DM	883	0	941	42	0
14	AN	774	0	827	52	0
14	DN	774	0	827	25	0
15	AO	716	0	739	33	0
15	DO	716	0	739	18	0
16	AP	649	0	666	20	0
16	DP	649	0	666	36	0
17	AQ	648	0	691	17	0
17	DQ	648	0	691	15	0
18	AR	455	0	478	20	0
18	DR	455	0	478	15	0
19	AS	637	0	665	38	0
19	DS	637	0	665	26	0
20	AT	665	0	714	15	0
20	DT	665	0	714	20	0
21	AU	425	0	449	24	0
21	DU	425	0	449	35	0
22	AV	1419	0	1465	45	0
23	AW	324	0	162	4	0
23	DV	346	0	173	15	0
24	AX	1623	0	821	20	0
24	DW	1623	0	821	8	0
25	BA	62195	0	31279	791	0
25	CA	62195	0	31280	774	0
26	BB	2549	0	1291	25	0
26	CB	2529	0	1281	20	0
27	BC	2082	0	2157	64	0
27	CC	2082	0	2157	61	0
28	BD	1565	0	1616	46	0
28	CD	1565	0	1616	31	0
29	BE	1552	0	1619	39	0
29	CE	1552	0	1619	28	0
30	BF	1410	0	1447	53	0
30	CF	1410	0	1447	49	0
31	BG	1323	0	1374	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
31	CG	1323	0	1374	22	0
32	BH	1107	0	1139	90	0
32	CH	1107	0	1138	62	1
33	BI	1032	0	1088	17	0
33	CI	1032	0	1088	17	0
34	BJ	1129	0	1162	24	0
34	CJ	1129	0	1162	27	0
35	BK	938	0	1012	31	0
35	CK	938	0	1012	34	0
36	BL	1045	0	1117	35	0
36	CL	1045	0	1117	39	0
37	BM	1074	0	1157	23	0
37	CM	1074	0	1157	24	0
38	BN	960	0	1000	26	0
38	CN	960	0	1000	25	0
39	BO	892	0	922	24	0
39	CO	892	0	923	32	0
40	BP	917	0	965	30	0
40	CP	917	0	965	31	0
41	BQ	947	0	1022	36	0
41	CQ	947	0	1022	35	0
42	BR	816	0	839	25	0
42	CR	816	0	839	24	0
43	BS	857	0	922	25	0
43	CS	857	0	922	22	0
44	BT	738	0	807	37	0
44	CT	738	0	806	31	0
45	BU	779	0	834	25	0
45	CU	779	0	834	32	0
46	BV	753	0	780	18	0
46	CV	753	0	780	8	0
47	BW	580	0	594	17	0
47	CW	569	0	581	11	0
48	BX	625	0	655	15	0
48	CX	625	0	655	16	0
49	BY	509	0	543	22	0
49	CY	509	0	543	20	0
50	BZ	449	0	491	6	0
50	CZ	449	0	491	11	0
51	B0	444	0	461	15	0
51	C0	444	0	461	10	0
52	B1	409	0	440	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
52	C1	409	0	440	5	0
53	B2	377	0	418	9	0
53	C2	377	0	418	8	0
54	B3	504	0	574	38	0
54	C3	504	0	574	11	0
55	B4	302	0	340	13	0
55	C4	302	0	340	4	0
56	B5	1142	0	865	13	0
57	AA	71	0	0	0	0
57	AN	1	0	0	0	0
57	BA	189	0	0	0	0
57	BB	4	0	0	0	0
57	BL	2	0	0	0	0
57	BO	1	0	0	0	0
57	BQ	1	0	0	0	0
57	CA	165	0	0	0	0
57	CB	3	0	0	0	0
57	CD	1	0	0	0	0
57	CQ	1	0	0	0	0
57	DA	53	0	0	0	0
57	DD	2	0	0	0	0
57	DN	1	0	0	0	0
58	AA	42	45	45	9	0
58	BA	210	180	225	33	0
58	CA	210	135	224	51	0
58	DA	84	90	90	13	0
59	B4	1	0	0	0	0
59	C4	1	0	0	0	0
60	AA	192	0	0	17	0
60	AE	3	0	0	0	0
60	AL	1	0	0	0	0
60	AN	2	0	0	1	0
60	AT	5	0	0	0	0
60	B3	1	0	0	0	0
60	B4	1	0	0	0	0
60	BA	626	0	0	92	0
60	BB	14	0	0	3	0
60	BC	7	0	0	0	0
60	BD	2	0	0	1	0
60	BE	1	0	0	0	0
60	BF	1	0	0	0	0
60	BL	8	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
60	BN	2	0	0	0	0
60	BQ	1	0	0	0	0
60	BS	1	0	0	0	0
60	BT	3	0	0	2	0
60	C3	1	0	0	0	0
60	C4	2	0	0	0	0
60	CA	608	0	0	82	0
60	CB	14	0	0	3	0
60	CC	10	0	0	1	0
60	CD	5	0	0	1	0
60	CE	3	0	0	0	0
60	CJ	2	0	0	2	0
60	CL	7	0	0	0	0
60	CN	2	0	0	0	0
60	CT	3	0	0	1	0
60	CU	1	0	0	0	0
60	DA	184	0	0	22	0
60	DD	2	0	0	1	0
60	DK	2	0	0	0	0
60	DL	2	0	0	0	0
60	DN	4	0	0	2	0
60	DT	3	0	0	0	0
60	DU	1	0	0	0	0
All	All	294034	450	196884	4955	1

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 10.

The worst 5 of 4955 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)
7:AG:103:TRP:CE2	7:AG:103:TRP:CZ2	1.76	1.63
7:AG:103:TRP:CH2	7:AG:103:TRP:CZ3	1.85	1.61
7:AG:100:ALA:HA	7:AG:103:TRP:CD2	1.52	1.44
7:AG:103:TRP:CZ3	7:AG:103:TRP:CE3	2.06	1.44
32:BH:124:THR:HG21	32:BH:128:HIS:NE2	1.32	1.44

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:358:U:OP1	32:CH:123:ARG:NH2[4_455]	1.91	0.29

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	
2	AB	216/241 (90%)	120 (56%)	54 (25%)	42 (19%)	0	0
2	DB	216/241 (90%)	130 (60%)	50 (23%)	36 (17%)	0	1
3	AC	204/233 (88%)	148 (72%)	34 (17%)	22 (11%)	0	2
3	DC	204/233 (88%)	154 (76%)	33 (16%)	17 (8%)	1	4
4	AD	203/206 (98%)	115 (57%)	55 (27%)	33 (16%)	0	1
4	DD	203/206 (98%)	111 (55%)	41 (20%)	51 (25%)	0	0
5	AE	148/167 (89%)	116 (78%)	19 (13%)	13 (9%)	1	4
5	DE	148/167 (89%)	106 (72%)	25 (17%)	17 (12%)	0	2
6	AF	98/135 (73%)	62 (63%)	22 (22%)	14 (14%)	0	1
6	DF	98/135 (73%)	74 (76%)	11 (11%)	13 (13%)	0	1
7	AG	149/179 (83%)	88 (59%)	43 (29%)	18 (12%)	0	2
7	DG	149/179 (83%)	112 (75%)	24 (16%)	13 (9%)	1	4
8	AH	127/130 (98%)	92 (72%)	27 (21%)	8 (6%)	1	9
8	DH	127/130 (98%)	97 (76%)	14 (11%)	16 (13%)	0	1
9	AI	125/130 (96%)	84 (67%)	28 (22%)	13 (10%)	0	3
9	DI	125/130 (96%)	83 (66%)	23 (18%)	19 (15%)	0	1
10	AJ	96/103 (93%)	60 (62%)	22 (23%)	14 (15%)	0	1
10	DJ	96/103 (93%)	71 (74%)	17 (18%)	8 (8%)	1	4
11	AK	115/129 (89%)	67 (58%)	30 (26%)	18 (16%)	0	1
11	DK	115/129 (89%)	85 (74%)	17 (15%)	13 (11%)	0	2
12	AL	121/124 (98%)	94 (78%)	14 (12%)	13 (11%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
12	DL	121/124 (98%)	72 (60%)	32 (26%)	17 (14%)	0	1
13	AM	112/118 (95%)	63 (56%)	29 (26%)	20 (18%)	0	0
13	DM	112/118 (95%)	71 (63%)	26 (23%)	15 (13%)	0	1
14	AN	92/101 (91%)	65 (71%)	18 (20%)	9 (10%)	1	3
14	DN	92/101 (91%)	63 (68%)	15 (16%)	14 (15%)	0	1
15	AO	86/89 (97%)	65 (76%)	17 (20%)	4 (5%)	3	16
15	DO	86/89 (97%)	66 (77%)	14 (16%)	6 (7%)	1	7
16	AP	80/82 (98%)	48 (60%)	26 (32%)	6 (8%)	1	6
16	DP	80/82 (98%)	48 (60%)	25 (31%)	7 (9%)	1	4
17	AQ	78/84 (93%)	55 (70%)	11 (14%)	12 (15%)	0	1
17	DQ	78/84 (93%)	47 (60%)	21 (27%)	10 (13%)	0	1
18	AR	53/75 (71%)	37 (70%)	9 (17%)	7 (13%)	0	1
18	DR	53/75 (71%)	38 (72%)	7 (13%)	8 (15%)	0	1
19	AS	77/92 (84%)	45 (58%)	17 (22%)	15 (20%)	0	0
19	DS	77/92 (84%)	53 (69%)	16 (21%)	8 (10%)	0	3
20	AT	83/87 (95%)	69 (83%)	6 (7%)	8 (10%)	1	3
20	DT	83/87 (95%)	62 (75%)	15 (18%)	6 (7%)	1	6
21	AU	49/71 (69%)	20 (41%)	16 (33%)	13 (26%)	0	0
21	DU	49/71 (69%)	25 (51%)	9 (18%)	15 (31%)	0	0
22	AV	181/185 (98%)	137 (76%)	30 (17%)	14 (8%)	1	5
27	BC	269/273 (98%)	216 (80%)	37 (14%)	16 (6%)	2	11
27	CC	269/273 (98%)	214 (80%)	41 (15%)	14 (5%)	2	14
28	BD	207/209 (99%)	176 (85%)	22 (11%)	9 (4%)	3	18
28	CD	207/209 (99%)	170 (82%)	25 (12%)	12 (6%)	2	11
29	BE	199/201 (99%)	160 (80%)	30 (15%)	9 (4%)	3	17
29	CE	199/201 (99%)	156 (78%)	31 (16%)	12 (6%)	2	10
30	BF	175/179 (98%)	127 (73%)	32 (18%)	16 (9%)	1	4
30	CF	175/179 (98%)	122 (70%)	37 (21%)	16 (9%)	1	4
31	BG	174/177 (98%)	133 (76%)	29 (17%)	12 (7%)	1	7
31	CG	174/177 (98%)	135 (78%)	27 (16%)	12 (7%)	1	7
32	BH	147/149 (99%)	84 (57%)	40 (27%)	23 (16%)	0	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
32	CH	147/149 (99%)	93 (63%)	36 (24%)	18 (12%)	0	1
33	BI	139/142 (98%)	77 (55%)	37 (27%)	25 (18%)	0	0
33	CI	139/142 (98%)	84 (60%)	39 (28%)	16 (12%)	0	2
34	BJ	140/142 (99%)	123 (88%)	11 (8%)	6 (4%)	3	18
34	CJ	140/142 (99%)	123 (88%)	13 (9%)	4 (3%)	5	28
35	BK	120/123 (98%)	99 (82%)	10 (8%)	11 (9%)	1	4
35	CK	120/123 (98%)	92 (77%)	22 (18%)	6 (5%)	2	15
36	BL	141/144 (98%)	114 (81%)	15 (11%)	12 (8%)	1	4
36	CL	141/144 (98%)	98 (70%)	29 (21%)	14 (10%)	1	3
37	BM	134/136 (98%)	110 (82%)	16 (12%)	8 (6%)	2	10
37	CM	134/136 (98%)	109 (81%)	17 (13%)	8 (6%)	2	10
38	BN	118/127 (93%)	98 (83%)	14 (12%)	6 (5%)	2	14
38	CN	118/127 (93%)	83 (70%)	32 (27%)	3 (2%)	6	32
39	BO	114/117 (97%)	90 (79%)	18 (16%)	6 (5%)	2	13
39	CO	114/117 (97%)	80 (70%)	23 (20%)	11 (10%)	1	3
40	BP	112/115 (97%)	94 (84%)	15 (13%)	3 (3%)	6	30
40	CP	112/115 (97%)	91 (81%)	16 (14%)	5 (4%)	3	17
41	BQ	115/118 (98%)	106 (92%)	8 (7%)	1 (1%)	20	62
41	CQ	115/118 (98%)	99 (86%)	13 (11%)	3 (3%)	6	31
42	BR	101/103 (98%)	89 (88%)	7 (7%)	5 (5%)	2	15
42	CR	101/103 (98%)	76 (75%)	18 (18%)	7 (7%)	1	7
43	BS	108/110 (98%)	96 (89%)	10 (9%)	2 (2%)	9	41
43	CS	108/110 (98%)	82 (76%)	16 (15%)	10 (9%)	1	4
44	BT	91/100 (91%)	72 (79%)	11 (12%)	8 (9%)	1	4
44	CT	91/100 (91%)	72 (79%)	11 (12%)	8 (9%)	1	4
45	BU	100/104 (96%)	79 (79%)	14 (14%)	7 (7%)	1	7
45	CU	100/104 (96%)	67 (67%)	23 (23%)	10 (10%)	1	3
46	BV	92/94 (98%)	81 (88%)	11 (12%)	0	100	100
46	CV	92/94 (98%)	74 (80%)	14 (15%)	4 (4%)	3	18
47	BW	74/85 (87%)	60 (81%)	9 (12%)	5 (7%)	1	7
47	CW	73/85 (86%)	60 (82%)	10 (14%)	3 (4%)	3	19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
48	BX	75/78 (96%)	62 (83%)	9 (12%)	4 (5%)	2	13
48	CX	75/78 (96%)	63 (84%)	10 (13%)	2 (3%)	6	30
49	BY	61/63 (97%)	46 (75%)	9 (15%)	6 (10%)	1	3
49	CY	61/63 (97%)	44 (72%)	12 (20%)	5 (8%)	1	5
50	BZ	56/59 (95%)	45 (80%)	11 (20%)	0	100	100
50	CZ	56/59 (95%)	47 (84%)	9 (16%)	0	100	100
51	B0	54/57 (95%)	45 (83%)	4 (7%)	5 (9%)	1	4
51	C0	54/57 (95%)	42 (78%)	4 (7%)	8 (15%)	0	1
52	B1	48/55 (87%)	31 (65%)	12 (25%)	5 (10%)	0	3
52	C1	48/55 (87%)	37 (77%)	11 (23%)	0	100	100
53	B2	44/46 (96%)	37 (84%)	6 (14%)	1 (2%)	7	35
53	C2	44/46 (96%)	38 (86%)	4 (9%)	2 (4%)	3	17
54	B3	62/65 (95%)	51 (82%)	7 (11%)	4 (6%)	1	8
54	C3	62/65 (95%)	52 (84%)	8 (13%)	2 (3%)	5	26
55	B4	36/38 (95%)	31 (86%)	3 (8%)	2 (6%)	2	12
55	C4	36/38 (95%)	30 (83%)	5 (14%)	1 (3%)	6	29
56	B5	183/228 (80%)	80 (44%)	66 (36%)	37 (20%)	0	0
All	All	11599/12383 (94%)	8463 (73%)	2041 (18%)	1095 (9%)	1	3

5 of 1095 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	AB	23	TRP
2	AB	31	ILE
2	AB	34	ALA
2	AB	52	GLU
2	AB	73	LYS

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	AB	180/199 (90%)	145 (81%)	35 (19%)	1	9
2	DB	180/199 (90%)	142 (79%)	38 (21%)	1	6
3	AC	170/190 (90%)	138 (81%)	32 (19%)	2	10
3	DC	170/190 (90%)	140 (82%)	30 (18%)	2	11
4	AD	172/173 (99%)	126 (73%)	46 (27%)	0	3
4	DD	172/173 (99%)	130 (76%)	42 (24%)	1	3
5	AE	113/126 (90%)	91 (80%)	22 (20%)	1	8
5	DE	113/126 (90%)	89 (79%)	24 (21%)	1	6
6	AF	87/116 (75%)	68 (78%)	19 (22%)	1	6
6	DF	87/116 (75%)	70 (80%)	17 (20%)	1	8
7	AG	124/147 (84%)	101 (82%)	23 (18%)	2	10
7	DG	124/147 (84%)	106 (86%)	18 (14%)	4	17
8	AH	104/105 (99%)	85 (82%)	19 (18%)	2	10
8	DH	104/105 (99%)	87 (84%)	17 (16%)	3	13
9	AI	105/107 (98%)	78 (74%)	27 (26%)	0	3
9	DI	105/107 (98%)	82 (78%)	23 (22%)	1	5
10	AJ	86/90 (96%)	66 (77%)	20 (23%)	1	4
10	DJ	86/90 (96%)	67 (78%)	19 (22%)	1	5
11	AK	90/99 (91%)	72 (80%)	18 (20%)	1	8
11	DK	90/99 (91%)	75 (83%)	15 (17%)	2	13
12	AL	103/104 (99%)	74 (72%)	29 (28%)	0	2
12	DL	103/104 (99%)	82 (80%)	21 (20%)	1	7
13	AM	92/96 (96%)	68 (74%)	24 (26%)	0	3
13	DM	92/96 (96%)	74 (80%)	18 (20%)	1	8
14	AN	79/84 (94%)	60 (76%)	19 (24%)	1	4
14	DN	79/84 (94%)	59 (75%)	20 (25%)	0	3
15	AO	76/77 (99%)	59 (78%)	17 (22%)	1	5
15	DO	76/77 (99%)	60 (79%)	16 (21%)	1	6
16	AP	65/65 (100%)	47 (72%)	18 (28%)	0	2
16	DP	65/65 (100%)	51 (78%)	14 (22%)	1	6
17	AQ	74/78 (95%)	61 (82%)	13 (18%)	2	11
17	DQ	74/78 (95%)	54 (73%)	20 (27%)	0	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
18	AR	48/65 (74%)	33 (69%)	15 (31%)	0	1
18	DR	48/65 (74%)	36 (75%)	12 (25%)	1	3
19	AS	70/79 (89%)	47 (67%)	23 (33%)	0	1
19	DS	70/79 (89%)	56 (80%)	14 (20%)	1	8
20	AT	65/66 (98%)	48 (74%)	17 (26%)	0	3
20	DT	65/66 (98%)	51 (78%)	14 (22%)	1	6
21	AU	44/61 (72%)	31 (70%)	13 (30%)	0	2
21	DU	44/61 (72%)	30 (68%)	14 (32%)	0	1
22	AV	157/160 (98%)	128 (82%)	29 (18%)	2	10
27	BC	216/218 (99%)	173 (80%)	43 (20%)	1	8
27	CC	216/218 (99%)	176 (82%)	40 (18%)	2	10
28	BD	164/164 (100%)	131 (80%)	33 (20%)	1	7
28	CD	164/164 (100%)	135 (82%)	29 (18%)	2	11
29	BE	165/165 (100%)	142 (86%)	23 (14%)	4	18
29	CE	165/165 (100%)	137 (83%)	28 (17%)	2	12
30	BF	148/150 (99%)	118 (80%)	30 (20%)	1	7
30	CF	148/150 (99%)	125 (84%)	23 (16%)	3	15
31	BG	137/138 (99%)	109 (80%)	28 (20%)	1	7
31	CG	137/138 (99%)	112 (82%)	25 (18%)	2	10
32	BH	113/114 (99%)	90 (80%)	23 (20%)	1	7
32	CH	113/114 (99%)	85 (75%)	28 (25%)	1	3
33	BI	109/110 (99%)	94 (86%)	15 (14%)	4	19
33	CI	109/110 (99%)	96 (88%)	13 (12%)	6	25
34	BJ	116/116 (100%)	98 (84%)	18 (16%)	3	15
34	CJ	116/116 (100%)	97 (84%)	19 (16%)	2	13
35	BK	103/104 (99%)	86 (84%)	17 (16%)	2	13
35	CK	103/104 (99%)	79 (77%)	24 (23%)	1	4
36	BL	102/103 (99%)	78 (76%)	24 (24%)	1	4
36	CL	102/103 (99%)	82 (80%)	20 (20%)	1	8
37	BM	109/109 (100%)	96 (88%)	13 (12%)	6	25
37	CM	109/109 (100%)	90 (83%)	19 (17%)	2	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
38	BN	100/103 (97%)	80 (80%)	20 (20%)	1	8
38	CN	100/103 (97%)	91 (91%)	9 (9%)	11	40
39	BO	86/87 (99%)	65 (76%)	21 (24%)	1	3
39	CO	86/87 (99%)	63 (73%)	23 (27%)	0	3
40	BP	99/100 (99%)	78 (79%)	21 (21%)	1	6
40	CP	99/100 (99%)	86 (87%)	13 (13%)	5	21
41	BQ	89/90 (99%)	77 (86%)	12 (14%)	4	19
41	CQ	89/90 (99%)	74 (83%)	15 (17%)	2	12
42	BR	84/84 (100%)	74 (88%)	10 (12%)	6	25
42	CR	84/84 (100%)	73 (87%)	11 (13%)	5	21
43	BS	93/93 (100%)	77 (83%)	16 (17%)	2	12
43	CS	93/93 (100%)	72 (77%)	21 (23%)	1	5
44	BT	80/84 (95%)	65 (81%)	15 (19%)	2	10
44	CT	80/84 (95%)	66 (82%)	14 (18%)	2	11
45	BU	83/85 (98%)	63 (76%)	20 (24%)	1	4
45	CU	83/85 (98%)	67 (81%)	16 (19%)	1	9
46	BV	78/78 (100%)	68 (87%)	10 (13%)	5	22
46	CV	78/78 (100%)	63 (81%)	15 (19%)	1	9
47	BW	57/63 (90%)	46 (81%)	11 (19%)	1	9
47	CW	56/63 (89%)	48 (86%)	8 (14%)	4	18
48	BX	67/68 (98%)	55 (82%)	12 (18%)	2	11
48	CX	67/68 (98%)	55 (82%)	12 (18%)	2	11
49	BY	55/55 (100%)	43 (78%)	12 (22%)	1	6
49	CY	55/55 (100%)	45 (82%)	10 (18%)	2	10
50	BZ	48/49 (98%)	40 (83%)	8 (17%)	2	13
50	CZ	48/49 (98%)	40 (83%)	8 (17%)	2	13
51	B0	47/48 (98%)	44 (94%)	3 (6%)	20	57
51	C0	47/48 (98%)	41 (87%)	6 (13%)	5	22
52	B1	45/49 (92%)	39 (87%)	6 (13%)	4	20
52	C1	45/49 (92%)	38 (84%)	7 (16%)	3	15
53	B2	38/38 (100%)	31 (82%)	7 (18%)	2	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
53	C2	38/38 (100%)	31 (82%)	7 (18%)	2	10
54	B3	51/52 (98%)	43 (84%)	8 (16%)	3	14
54	C3	51/52 (98%)	42 (82%)	9 (18%)	2	11
55	B4	34/34 (100%)	26 (76%)	8 (24%)	1	4
55	C4	34/34 (100%)	27 (79%)	7 (21%)	1	7
56	B5	61/180 (34%)	52 (85%)	9 (15%)	3	16
All	All	9543/10096 (94%)	7684 (80%)	1859 (20%)	1	8

5 of 1859 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
43	BS	97	LEU
28	CD	177	VAL
13	DM	14	HIS
45	BU	34	ILE
53	B2	15	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 62 such sidechains are listed below:

Mol	Chain	Res	Type
34	BJ	47	HIS
56	B5	66	HIS
12	DL	77	HIS
45	BU	44	HIS
27	CC	116	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	1537/1542 (99%)	349 (22%)	15 (0%)
1	DA	1538/1542 (99%)	330 (21%)	17 (1%)
23	AW	14/16 (87%)	5 (35%)	0
23	DV	15/16 (93%)	6 (40%)	0
24	AX	75/76 (98%)	27 (36%)	6 (8%)
24	DW	75/76 (98%)	20 (26%)	0
25	BA	2896/2904 (99%)	590 (20%)	35 (1%)
25	CA	2895/2904 (99%)	618 (21%)	34 (1%)
26	BB	118/120 (98%)	16 (13%)	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
26	CB	117/120 (97%)	27 (23%)	0
All	All	9280/9316 (99%)	1988 (21%)	107 (1%)

5 of 1988 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	5	U
1	AA	6	G
1	AA	9	G
1	AA	28	A
1	AA	30	U

5 of 107 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
25	BA	2324	U
25	CA	479	A
1	DA	723	U
25	BA	2326	C
25	BA	2756	U

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 510 ligands modelled in this entry, 497 are monoatomic - leaving 13 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
58	PAR	AA	1672	-	45,45,45	0.76	0	60,67,67	1.18	4 (6%)
58	PAR	BA	3001	-	45,45,45	0.72	0	60,67,67	1.07	5 (8%)
58	PAR	BA	3002	-	45,45,45	0.79	1 (2%)	60,67,67	1.03	4 (6%)
58	PAR	BA	3003	-	45,45,45	0.73	0	60,67,67	1.03	4 (6%)
58	PAR	BA	3004	-	45,45,45	0.71	0	60,67,67	1.08	5 (8%)
58	PAR	BA	3005	-	45,45,45	0.55	0	60,67,67	1.15	4 (6%)
58	PAR	CA	3166	-	45,45,45	0.49	0	60,67,67	0.98	4 (6%)
58	PAR	CA	3167	-	45,45,45	0.58	0	60,67,67	0.92	3 (5%)
58	PAR	CA	3168	-	45,45,45	0.52	0	60,67,67	1.38	6 (10%)
58	PAR	CA	3169	-	45,45,45	0.49	0	60,67,67	0.99	4 (6%)
58	PAR	CA	3170	-	45,45,45	0.50	0	60,67,67	0.96	5 (8%)
58	PAR	DA	1654	-	45,45,45	0.53	0	60,67,67	0.85	2 (3%)
58	PAR	DA	1655	-	45,45,45	0.50	0	60,67,67	0.95	3 (5%)

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
58	PAR	AA	1672	-	-	0/18/94/94	0/4/4/4
58	PAR	BA	3001	-	-	0/18/94/94	0/4/4/4
58	PAR	BA	3002	-	-	0/18/94/94	0/4/4/4
58	PAR	BA	3003	-	-	0/18/94/94	0/4/4/4
58	PAR	BA	3004	-	-	0/18/94/94	0/4/4/4
58	PAR	BA	3005	-	-	0/18/94/94	0/4/4/4
58	PAR	CA	3166	-	-	0/18/94/94	0/4/4/4
58	PAR	CA	3167	-	-	0/18/94/94	0/4/4/4
58	PAR	CA	3168	-	-	0/18/94/94	0/4/4/4
58	PAR	CA	3169	-	-	0/18/94/94	0/4/4/4
58	PAR	CA	3170	-	-	0/18/94/94	0/4/4/4
58	PAR	DA	1654	-	-	0/18/94/94	0/4/4/4
58	PAR	DA	1655	-	-	0/18/94/94	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
58	BA	3002	PAR	C13-C23	-2.00	1.50	1.52

The worst 5 of 53 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
58	CA	3168	PAR	C14-O33-C33	-4.61	106.77	118.00
58	CA	3168	PAR	C11-O11-C42	-4.58	106.83	118.00
58	CA	3168	PAR	C13-O52-C52	-4.56	106.88	118.00
58	BA	3004	PAR	O52-C13-O43	-3.78	107.33	111.43
58	BA	3005	PAR	C13-O52-C52	-3.68	109.03	118.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

13 monomers are involved in 106 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
58	AA	1672	PAR	9	0
58	BA	3001	PAR	3	0
58	BA	3002	PAR	2	0
58	BA	3003	PAR	4	0
58	BA	3004	PAR	3	0
58	BA	3005	PAR	21	0
58	CA	3166	PAR	14	0
58	CA	3167	PAR	7	0
58	CA	3168	PAR	5	0
58	CA	3169	PAR	20	0
58	CA	3170	PAR	5	0
58	DA	1654	PAR	5	0
58	DA	1655	PAR	8	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	AA	1538/1542 (99%)	-0.65	2 (0%) 95 88	44, 86, 166, 303	0
1	DA	1539/1542 (99%)	-0.63	1 (0%) 95 88	47, 85, 145, 285	0
2	AB	218/241 (90%)	0.81	28 (12%) 4 2	64, 133, 196, 238	0
2	DB	218/241 (90%)	0.48	14 (6%) 20 7	68, 115, 176, 210	0
3	AC	206/233 (88%)	-0.44	0 100 100	44, 77, 114, 170	0
3	DC	206/233 (88%)	-0.17	0 100 100	49, 90, 145, 213	0
4	AD	205/206 (99%)	0.03	7 (3%) 46 20	62, 104, 170, 243	0
4	DD	205/206 (99%)	0.01	9 (4%) 35 14	47, 92, 166, 215	0
5	AE	150/167 (89%)	-0.16	1 (0%) 87 67	53, 81, 124, 222	0
5	DE	150/167 (89%)	-0.39	0 100 100	44, 79, 132, 171	0
6	AF	100/135 (74%)	-0.12	1 (1%) 82 58	67, 122, 177, 211	0
6	DF	100/135 (74%)	-0.27	2 (2%) 65 36	50, 94, 135, 153	0
7	AG	151/179 (84%)	0.18	7 (4%) 33 13	62, 119, 169, 239	0
7	DG	151/179 (84%)	-0.05	4 (2%) 56 27	61, 104, 150, 211	0
8	AH	129/130 (99%)	-0.05	3 (2%) 61 31	59, 92, 136, 159	0
8	DH	129/130 (99%)	-0.20	2 (1%) 72 44	49, 84, 119, 174	0
9	AI	127/130 (97%)	0.33	9 (7%) 17 6	57, 108, 158, 215	0
9	DI	127/130 (97%)	0.23	5 (3%) 40 16	70, 116, 174, 197	0
10	AJ	98/103 (95%)	0.47	9 (9%) 10 4	57, 111, 175, 216	0
10	DJ	98/103 (95%)	-0.03	3 (3%) 49 22	65, 111, 168, 200	0
11	AK	117/129 (90%)	0.36	12 (10%) 7 3	58, 108, 177, 232	0
11	DK	117/129 (90%)	-0.20	4 (3%) 46 20	47, 75, 131, 186	0
12	AL	123/124 (99%)	-0.23	3 (2%) 59 30	33, 65, 115, 216	0
12	DL	123/124 (99%)	0.05	1 (0%) 86 64	33, 71, 115, 205	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AM	114/118 (96%)	0.70	11 (9%) 9 3	73, 132, 191, 263	0
13	DM	114/118 (96%)	0.35	9 (7%) 13 5	92, 135, 174, 254	0
14	AN	96/101 (95%)	0.05	4 (4%) 37 15	59, 95, 188, 237	0
14	DN	96/101 (95%)	0.05	2 (2%) 64 34	62, 104, 164, 215	0
15	AO	88/89 (98%)	0.10	3 (3%) 46 20	60, 102, 166, 216	0
15	DO	88/89 (98%)	-0.10	2 (2%) 61 31	55, 85, 131, 156	0
16	AP	82/82 (100%)	0.40	6 (7%) 16 6	61, 86, 144, 198	0
16	DP	82/82 (100%)	0.43	6 (7%) 16 6	53, 79, 141, 227	0
17	AQ	80/84 (95%)	0.29	5 (6%) 21 7	52, 97, 152, 198	0
17	DQ	80/84 (95%)	0.51	6 (7%) 15 6	58, 96, 161, 233	0
18	AR	55/75 (73%)	0.01	2 (3%) 43 18	57, 97, 151, 178	0
18	DR	55/75 (73%)	-0.25	2 (3%) 43 18	57, 83, 134, 176	0
19	AS	79/92 (85%)	0.76	8 (10%) 8 3	91, 129, 184, 223	0
19	DS	79/92 (85%)	0.79	9 (11%) 6 2	83, 129, 181, 234	0
20	AT	85/87 (97%)	0.25	3 (3%) 44 19	54, 92, 134, 194	0
20	DT	85/87 (97%)	0.15	2 (2%) 59 30	64, 96, 132, 179	0
21	AU	51/71 (71%)	0.09	0 100 100	67, 117, 167, 203	0
21	DU	51/71 (71%)	-0.17	2 (3%) 40 16	54, 105, 159, 224	0
22	AV	183/185 (98%)	-0.12	3 (1%) 72 44	20, 86, 171, 232	0
23	AW	15/16 (93%)	-0.19	0 100 100	57, 128, 176, 205	0
23	DV	16/16 (100%)	-0.30	1 (6%) 21 7	44, 99, 143, 191	0
24	AX	76/76 (100%)	0.39	7 (9%) 10 4	44, 167, 277, 328	0
24	DW	76/76 (100%)	-0.48	0 100 100	51, 93, 135, 164	0
25	BA	2897/2904 (99%)	-0.44	95 (3%) 47 21	22, 52, 194, 368	0
25	CA	2897/2904 (99%)	-0.51	46 (1%) 72 44	37, 75, 205, 328	0
26	BB	119/120 (99%)	-0.82	0 100 100	41, 71, 101, 165	0
26	CB	118/120 (98%)	-0.81	0 100 100	63, 118, 153, 186	0
27	BC	271/273 (99%)	-0.47	0 100 100	32, 60, 94, 149	0
27	CC	271/273 (99%)	-0.34	1 (0%) 92 77	38, 73, 106, 185	0
28	BD	209/209 (100%)	-0.53	0 100 100	21, 45, 77, 170	0
28	CD	209/209 (100%)	-0.37	0 100 100	41, 68, 109, 171	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
29	BE	201/201 (100%)	-0.34	0	100	100	23, 56, 105, 161	0
29	CE	201/201 (100%)	-0.18	0	100	100	32, 79, 125, 175	0
30	BF	177/179 (98%)	0.48	15 (8%)	11	4	44, 104, 177, 231	0
30	CF	177/179 (98%)	0.90	32 (18%)	1	1	87, 140, 202, 243	0
31	BG	176/177 (99%)	-0.14	3 (1%)	70	42	32, 72, 114, 158	0
31	CG	176/177 (99%)	0.40	7 (3%)	39	16	72, 112, 153, 195	0
32	BH	149/149 (100%)	0.24	7 (4%)	32	13	28, 118, 177, 265	0
32	CH	149/149 (100%)	0.57	14 (9%)	9	3	28, 133, 188, 220	0
33	BI	141/142 (99%)	2.88	81 (57%)	0	0	118, 197, 266, 320	0
33	CI	141/142 (99%)	3.45	91 (64%)	0	0	137, 210, 253, 292	0
34	BJ	142/142 (100%)	-0.57	0	100	100	21, 44, 79, 102	0
34	CJ	142/142 (100%)	-0.31	0	100	100	37, 65, 99, 167	0
35	BK	122/123 (99%)	-0.59	0	100	100	23, 49, 77, 104	0
35	CK	122/123 (99%)	-0.25	0	100	100	42, 69, 112, 148	0
36	BL	143/144 (99%)	-0.52	0	100	100	24, 54, 83, 118	0
36	CL	143/144 (99%)	-0.15	2 (1%)	75	49	39, 82, 125, 180	0
37	BM	136/136 (100%)	-0.48	0	100	100	27, 51, 89, 121	0
37	CM	136/136 (100%)	0.20	4 (2%)	52	24	44, 78, 106, 141	0
38	BN	120/127 (94%)	-0.58	0	100	100	22, 47, 71, 156	0
38	CN	120/127 (94%)	-0.20	1 (0%)	86	64	46, 76, 103, 191	0
39	BO	116/117 (99%)	-0.18	0	100	100	37, 67, 99, 123	0
39	CO	116/117 (99%)	0.48	8 (6%)	18	6	70, 122, 164, 201	0
40	BP	114/115 (99%)	-0.53	2 (1%)	69	40	30, 56, 99, 206	0
40	CP	114/115 (99%)	-0.32	0	100	100	45, 77, 116, 140	0
41	BQ	117/118 (99%)	-0.63	0	100	100	15, 39, 71, 91	0
41	CQ	117/118 (99%)	-0.42	0	100	100	32, 61, 85, 114	0
42	BR	103/103 (100%)	-0.48	0	100	100	21, 52, 87, 118	0
42	CR	103/103 (100%)	-0.37	1 (0%)	82	58	32, 73, 108, 171	0
43	BS	110/110 (100%)	-0.47	0	100	100	20, 44, 76, 117	0
43	CS	110/110 (100%)	-0.37	0	100	100	31, 65, 103, 140	0
44	BT	93/100 (93%)	-0.10	1 (1%)	80	55	43, 60, 126, 213	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	CT	93/100 (93%)	0.06	2 (2%) 62 33	56, 94, 142, 202	0
45	BU	102/104 (98%)	-0.41	0 100 100	33, 61, 118, 185	0
45	CU	102/104 (98%)	0.42	9 (8%) 11 4	56, 90, 177, 215	0
46	BV	94/94 (100%)	-0.29	1 (1%) 80 55	41, 71, 109, 127	0
46	CV	94/94 (100%)	0.11	1 (1%) 80 55	66, 110, 151, 196	0
47	BW	76/85 (89%)	-0.41	1 (1%) 77 51	32, 54, 92, 126	0
47	CW	75/85 (88%)	0.14	1 (1%) 77 51	46, 82, 120, 146	0
48	BX	77/78 (98%)	-0.31	0 100 100	39, 59, 101, 111	0
48	CX	77/78 (98%)	-0.08	2 (2%) 56 27	38, 82, 121, 150	0
49	BY	63/63 (100%)	-0.17	3 (4%) 31 12	39, 69, 114, 179	0
49	CY	63/63 (100%)	0.29	3 (4%) 31 12	66, 103, 165, 221	0
50	BZ	58/59 (98%)	-0.33	0 100 100	30, 47, 75, 106	0
50	CZ	58/59 (98%)	-0.21	1 (1%) 70 42	45, 73, 112, 188	0
51	B0	56/57 (98%)	-0.62	0 100 100	22, 46, 94, 143	0
51	C0	56/57 (98%)	-0.33	1 (1%) 69 40	37, 73, 117, 194	0
52	B1	50/55 (90%)	2.19	17 (34%) 0 0	76, 105, 185, 196	0
52	C1	50/55 (90%)	1.03	7 (14%) 3 1	79, 109, 138, 164	0
53	B2	46/46 (100%)	-0.38	1 (2%) 62 33	27, 45, 76, 155	0
53	C2	46/46 (100%)	-0.23	0 100 100	39, 68, 92, 151	0
54	B3	64/65 (98%)	-0.42	0 100 100	27, 48, 74, 122	0
54	C3	64/65 (98%)	-0.29	0 100 100	38, 69, 96, 109	0
55	B4	38/38 (100%)	-0.32	0 100 100	29, 52, 91, 122	0
55	C4	38/38 (100%)	0.01	0 100 100	47, 83, 109, 119	0
56	B5	191/228 (83%)	5.26	156 (81%) 0 0	28, 219, 264, 311	0
All	All	21100/21699 (97%)	-0.17	837 (3%) 39 16	15, 80, 183, 368	0

The worst 5 of 837 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
56	B5	222	VAL	25.3
33	CI	1	ALA	21.3
56	B5	38	ASP	20.6
56	B5	223	ARG	19.9
56	B5	105	ASP	19.3

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
57	MG	CA	3019	1/1	0.95	0.99	54.82	252,252,252,252	0
57	MG	AA	1636	1/1	0.86	1.03	47.53	313,313,313,313	0
57	MG	CD	301	1/1	0.92	0.54	42.61	234,234,234,234	0
57	MG	CA	3098	1/1	0.97	0.64	36.84	242,242,242,242	0
57	MG	BA	3154	1/1	0.94	0.76	35.70	56,56,56,56	0
57	MG	DA	1633	1/1	0.95	0.85	25.00	264,264,264,264	0
58	PAR	DA	1655	42/42	0.83	0.50	24.54	10,10,10,10	0
57	MG	DA	1638	1/1	0.90	0.30	21.61	93,93,93,93	0
57	MG	AA	1645	1/1	0.89	0.32	19.91	75,75,75,75	0
57	MG	CA	3013	1/1	0.79	0.86	19.00	362,362,362,362	0
57	MG	BA	3067	1/1	0.97	0.31	17.90	218,218,218,218	0
57	MG	BA	3179	1/1	0.78	0.39	17.04	88,88,88,88	0
57	MG	BA	3134	1/1	0.99	0.28	14.95	143,143,143,143	0
58	PAR	CA	3168	42/42	0.73	0.61	14.76	10,10,10,10	0
57	MG	CA	3161	1/1	0.96	0.36	14.59	39,39,39,39	0
57	MG	BA	3109	1/1	0.75	0.37	13.99	189,189,189,189	0
57	MG	BA	3176	1/1	0.97	0.31	13.85	51,51,51,51	0
57	MG	CA	3094	1/1	0.95	0.20	11.17	94,94,94,94	0
58	PAR	CA	3167	42/42	0.89	0.36	10.75	10,10,10,10	0
57	MG	BA	3105	1/1	0.96	0.26	10.74	80,80,80,80	0
58	PAR	BA	3005	42/42	0.85	0.28	10.00	41,100,141,152	42
57	MG	BA	3157	1/1	0.98	0.30	9.89	44,44,44,44	0
57	MG	BA	3166	1/1	0.99	0.23	9.72	34,34,34,34	0
58	PAR	BA	3004	42/42	0.93	0.27	9.67	10,10,10,10	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	BA	3025	1/1	0.95	0.30	9.58	134,134,134,134	0
57	MG	CA	3064	1/1	0.94	0.24	9.57	118,118,118,118	0
57	MG	CA	3145	1/1	0.96	0.26	9.11	59,59,59,59	0
58	PAR	CA	3169	42/42	0.64	0.44	9.07	21,68,103,144	42
57	MG	DA	1644	1/1	0.98	0.29	8.63	62,62,62,62	0
58	PAR	BA	3003	42/42	0.89	0.39	8.40	10,10,10,10	0
57	MG	CA	3151	1/1	0.93	0.31	8.35	44,44,44,44	0
57	MG	AA	1669	1/1	0.96	0.43	7.98	58,58,58,58	0
57	MG	BA	3120	1/1	0.98	0.25	7.68	91,91,91,91	0
57	MG	CA	3142	1/1	0.99	0.21	7.28	33,33,33,33	0
58	PAR	BA	3001	42/42	0.93	0.30	6.77	10,10,10,10	0
57	MG	CA	3104	1/1	0.98	0.25	6.63	57,57,57,57	0
57	MG	AA	1631	1/1	0.97	0.23	6.40	115,115,115,115	0
57	MG	BA	3021	1/1	0.77	0.39	6.29	276,276,276,276	0
57	MG	BA	3014	1/1	0.96	0.20	6.23	101,101,101,101	0
57	MG	BA	3128	1/1	0.97	0.21	6.19	99,99,99,99	0
58	PAR	AA	1672	42/42	0.94	0.24	5.76	10,10,10,10	0
57	MG	BA	3044	1/1	0.91	0.25	5.52	62,62,62,62	0
57	MG	BA	3138	1/1	0.88	0.44	5.37	227,227,227,227	0
57	MG	BA	3077	1/1	0.97	0.22	5.10	63,63,63,63	0
58	PAR	BA	3002	42/42	0.95	0.23	4.76	10,10,10,10	0
58	PAR	CA	3166	42/42	0.83	0.32	4.76	38,86,109,128	42
57	MG	DA	1605	1/1	0.98	0.14	4.53	96,96,96,96	0
57	MG	BA	3155	1/1	0.99	0.26	4.51	23,23,23,23	0
57	MG	CA	3143	1/1	0.97	0.24	4.33	69,69,69,69	0
57	MG	AA	1614	1/1	0.97	0.21	4.25	146,146,146,146	0
59	ZN	B4	9501	1/1	0.98	0.54	4.16	386,386,386,386	0
57	MG	BA	3066	1/1	0.98	0.17	4.00	77,77,77,77	0
57	MG	BA	3112	1/1	0.96	0.21	3.97	29,29,29,29	0
57	MG	BA	3038	1/1	0.99	0.25	3.79	73,73,73,73	0
57	MG	CA	3116	1/1	0.98	0.19	3.75	94,94,94,94	0
57	MG	BA	3161	1/1	0.99	0.18	3.60	45,45,45,45	0
57	MG	AA	1619	1/1	0.91	0.19	3.32	70,70,70,70	0
57	MG	BA	3118	1/1	0.97	0.37	3.23	164,164,164,164	0
58	PAR	DA	1654	42/42	0.93	0.31	3.09	10,10,10,10	0
57	MG	BA	3045	1/1	0.96	0.40	3.06	217,217,217,217	0
57	MG	CA	3041	1/1	0.99	0.19	2.79	36,36,36,36	0
57	MG	BA	3188	1/1	0.99	0.15	2.73	39,39,39,39	0
57	MG	AA	1640	1/1	0.99	0.17	2.55	21,21,21,21	0
57	MG	CA	3133	1/1	0.96	0.17	2.43	182,182,182,182	0
57	MG	CA	3109	1/1	0.93	0.20	2.42	74,74,74,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3108	1/1	0.97	0.18	2.28	101,101,101,101	0
57	MG	AA	1612	1/1	0.99	0.13	2.19	94,94,94,94	0
57	MG	AA	1603	1/1	0.98	0.16	2.07	45,45,45,45	0
57	MG	CA	3048	1/1	0.92	0.19	2.02	157,157,157,157	0
57	MG	CA	3032	1/1	0.97	0.17	1.86	97,97,97,97	0
57	MG	CA	3106	1/1	0.97	0.20	1.82	33,33,33,33	0
57	MG	AA	1623	1/1	0.97	0.16	1.77	67,67,67,67	0
57	MG	CA	3150	1/1	0.97	0.24	1.77	54,54,54,54	0
57	MG	CA	3112	1/1	0.99	0.20	1.71	137,137,137,137	0
58	PAR	CA	3170	42/42	0.91	0.35	1.47	10,10,10,10	0
57	MG	AA	1635	1/1	0.98	0.15	1.46	234,234,234,234	0
57	MG	BA	3187	1/1	0.99	0.23	1.33	29,29,29,29	0
57	MG	BA	3075	1/1	0.92	0.17	1.29	88,88,88,88	0
57	MG	BA	3108	1/1	0.96	0.18	1.25	81,81,81,81	0
57	MG	BA	3184	1/1	0.98	0.22	1.04	40,40,40,40	0
57	MG	BA	3028	1/1	0.98	0.14	1.03	35,35,35,35	0
57	MG	BA	3114	1/1	0.97	0.15	0.95	65,65,65,65	0
57	MG	BA	3149	1/1	0.96	0.14	0.93	33,33,33,33	0
57	MG	BA	3083	1/1	0.94	0.14	0.74	84,84,84,84	0
57	MG	BA	3094	1/1	0.97	0.14	0.62	153,153,153,153	0
57	MG	BA	3147	1/1	0.98	0.15	0.59	54,54,54,54	0
57	MG	DA	1612	1/1	0.96	0.12	0.49	123,123,123,123	0
57	MG	CA	3008	1/1	0.94	0.17	0.33	74,74,74,74	0
57	MG	CA	3080	1/1	0.92	0.14	0.08	102,102,102,102	0
57	MG	BA	3163	1/1	0.98	0.18	-0.00	27,27,27,27	0
57	MG	BA	3027	1/1	0.97	0.16	-0.05	57,57,57,57	0
57	MG	CA	3074	1/1	0.91	0.14	-0.25	53,53,53,53	0
57	MG	BA	3111	1/1	0.98	0.14	-0.26	52,52,52,52	0
57	MG	DA	1607	1/1	0.98	0.15	-0.38	63,63,63,63	0
59	ZN	C4	9501	1/1	0.99	0.15	-0.38	109,109,109,109	0
57	MG	BA	3026	1/1	0.98	0.15	-0.39	47,47,47,47	0
57	MG	BA	3113	1/1	0.99	0.15	-0.39	96,96,96,96	0
57	MG	CA	3131	1/1	0.89	0.14	-0.42	75,75,75,75	0
57	MG	CA	3107	1/1	0.97	0.13	-0.42	95,95,95,95	0
57	MG	CA	3124	1/1	0.95	0.14	-0.44	125,125,125,125	0
57	MG	DA	1630	1/1	0.97	0.14	-0.48	72,72,72,72	0
57	MG	CA	3037	1/1	0.96	0.14	-0.53	90,90,90,90	0
57	MG	CA	3101	1/1	0.97	0.15	-0.57	58,58,58,58	0
57	MG	CA	3156	1/1	0.94	0.17	-0.63	45,45,45,45	0
57	MG	DA	1616	1/1	0.92	0.12	-0.73	98,98,98,98	0
57	MG	BB	3201	1/1	0.93	0.10	-0.75	107,107,107,107	0
57	MG	AA	1601	1/1	0.97	0.15	-0.76	50,50,50,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	DA	1632	1/1	0.93	0.14	-0.81	71,71,71,71	0
57	MG	CB	3201	1/1	0.94	0.11	-0.89	216,216,216,216	0
57	MG	BA	3101	1/1	0.97	0.16	-0.90	70,70,70,70	0
57	MG	AA	1604	1/1	0.72	0.09	-0.96	173,173,173,173	0
57	MG	DA	1618	1/1	0.97	0.14	-1.11	51,51,51,51	0
57	MG	AA	1617	1/1	0.97	0.10	-1.14	104,104,104,104	0
57	MG	BA	3124	1/1	0.99	0.13	-1.14	45,45,45,45	0
57	MG	CA	3024	1/1	0.97	0.14	-1.17	65,65,65,65	0
57	MG	CB	3202	1/1	0.96	0.08	-1.20	114,114,114,114	0
57	MG	BL	201	1/1	0.99	0.12	-1.22	21,21,21,21	0
57	MG	AA	1630	1/1	0.92	0.10	-1.23	113,113,113,113	0
57	MG	BA	3034	1/1	0.95	0.15	-1.28	41,41,41,41	0
57	MG	CA	3025	1/1	0.99	0.14	-1.29	121,121,121,121	0
57	MG	BA	3042	1/1	0.99	0.14	-1.35	32,32,32,32	0
57	MG	BA	3013	1/1	0.95	0.10	-1.39	40,40,40,40	0
57	MG	CA	3067	1/1	0.99	0.11	-1.39	64,64,64,64	0
57	MG	CA	3134	1/1	0.96	0.14	-1.41	55,55,55,55	0
57	MG	BA	3022	1/1	0.99	0.09	-1.44	28,28,28,28	0
57	MG	DA	1603	1/1	0.98	0.12	-1.57	37,37,37,37	0
57	MG	DA	1620	1/1	0.95	0.11	-1.75	78,78,78,78	0
57	MG	BA	3136	1/1	0.98	0.10	-1.77	61,61,61,61	0
57	MG	CA	3120	1/1	0.99	0.12	-1.86	62,62,62,62	0
57	MG	CA	3054	1/1	0.91	0.12	-1.94	41,41,41,41	0
57	MG	CA	3152	1/1	0.98	0.10	-2.00	44,44,44,44	0
57	MG	BA	3068	1/1	0.97	0.09	-2.13	27,27,27,27	0
57	MG	DA	1601	1/1	0.96	0.13	-2.15	47,47,47,47	0
57	MG	CA	3039	1/1	0.98	0.14	-2.15	58,58,58,58	0
57	MG	BA	3053	1/1	0.99	0.14	-2.15	29,29,29,29	0
57	MG	CA	3033	1/1	0.97	0.13	-2.15	56,56,56,56	0
57	MG	CA	3105	1/1	0.99	0.13	-2.20	44,44,44,44	0
57	MG	BA	3017	1/1	0.93	0.12	-2.21	56,56,56,56	0
57	MG	DA	1629	1/1	0.97	0.09	-2.48	86,86,86,86	0
57	MG	DA	1625	1/1	0.98	0.11	-2.51	42,42,42,42	0
57	MG	CA	3110	1/1	0.81	0.09	-2.55	61,61,61,61	0
57	MG	BA	3125	1/1	0.98	0.11	-2.64	36,36,36,36	0
57	MG	CA	3136	1/1	0.95	0.05	-2.75	107,107,107,107	0
57	MG	CA	3012	1/1	0.96	0.10	-2.86	30,30,30,30	0
57	MG	CA	3018	1/1	0.99	0.05	-2.86	41,41,41,41	0
57	MG	CA	3130	1/1	0.99	0.08	-3.02	44,44,44,44	0
57	MG	CA	3049	1/1	0.82	0.12	-3.11	76,76,76,76	0
57	MG	DA	1610	1/1	0.99	0.09	-3.15	67,67,67,67	0
57	MG	AA	1621	1/1	0.85	0.08	-3.28	157,157,157,157	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3051	1/1	1.00	0.07	-3.31	24,24,24,24	0
57	MG	CA	3132	1/1	0.96	0.09	-3.37	67,67,67,67	0
57	MG	BA	3057	1/1	0.98	0.12	-3.43	26,26,26,26	0
57	MG	BA	3062	1/1	0.98	0.10	-3.52	73,73,73,73	0
57	MG	BA	3140	1/1	0.97	0.06	-3.54	73,73,73,73	0
57	MG	DA	1614	1/1	0.85	0.06	-3.54	96,96,96,96	0
57	MG	CA	3059	1/1	0.97	0.10	-3.55	88,88,88,88	0
57	MG	BA	3032	1/1	0.96	0.11	-3.60	33,33,33,33	0
57	MG	BA	3010	1/1	0.89	0.11	-3.63	56,56,56,56	0
57	MG	BA	3069	1/1	0.99	0.09	-3.66	20,20,20,20	0
57	MG	BA	3031	1/1	0.96	0.08	-3.81	60,60,60,60	0
57	MG	CA	3121	1/1	0.99	0.06	-3.89	27,27,27,27	0
57	MG	BA	3054	1/1	0.95	0.07	-3.94	46,46,46,46	0
57	MG	AA	1610	1/1	0.99	0.04	-3.99	66,66,66,66	0
57	MG	BA	3070	1/1	0.98	0.12	-4.03	50,50,50,50	0
57	MG	CA	3060	1/1	0.97	0.08	-4.11	67,67,67,67	0
57	MG	CA	3023	1/1	0.99	0.10	-4.37	37,37,37,37	0
57	MG	CA	3072	1/1	0.96	0.11	-4.48	99,99,99,99	0
57	MG	CA	3050	1/1	0.99	0.09	-5.14	42,42,42,42	0
57	MG	CA	3063	1/1	0.97	0.09	-5.31	50,50,50,50	0
57	MG	AA	1622	1/1	0.97	0.06	-5.40	76,76,76,76	0
57	MG	BA	3135	1/1	0.97	0.07	-5.42	79,79,79,79	0
57	MG	CA	3129	1/1	0.99	0.09	-5.82	33,33,33,33	0
57	MG	CA	3097	1/1	0.93	0.07	-6.05	63,63,63,63	0
57	MG	CA	3118	1/1	0.98	0.05	-6.18	53,53,53,53	0
57	MG	CA	3002	1/1	0.99	0.04	-6.41	49,49,49,49	0
57	MG	CA	3005	1/1	0.96	0.05	-6.50	90,90,90,90	0
57	MG	BA	3072	1/1	0.98	0.08	-6.53	41,41,41,41	0
57	MG	CA	3017	1/1	0.99	0.10	-6.92	36,36,36,36	0
57	MG	CA	3066	1/1	0.98	0.06	-7.07	45,45,45,45	0
57	MG	AA	1626	1/1	0.98	0.09	-7.08	66,66,66,66	0
57	MG	BA	3133	1/1	0.99	0.11	-7.35	39,39,39,39	0
57	MG	BA	3007	1/1	0.99	0.07	-8.49	61,61,61,61	0
57	MG	BA	3183	1/1	0.94	0.17	-	37,37,37,37	0
57	MG	DA	1621	1/1	0.98	0.09	-	47,47,47,47	0
57	MG	BA	3047	1/1	0.99	0.08	-	42,42,42,42	0
57	MG	DA	1613	1/1	0.96	0.12	-	89,89,89,89	0
57	MG	CA	3096	1/1	0.96	0.07	-	100,100,100,100	0
57	MG	CA	3015	1/1	0.59	0.11	-	126,126,126,126	0
57	MG	BA	3089	1/1	0.96	0.15	-	44,44,44,44	0
57	MG	AA	1602	1/1	0.95	0.07	-	107,107,107,107	0
57	MG	AA	1624	1/1	0.97	0.07	-	99,99,99,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	AA	1667	1/1	0.85	0.27	-	81,81,81,81	0
57	MG	BA	3095	1/1	0.82	0.16	-	187,187,187,187	0
57	MG	CA	3083	1/1	0.91	0.14	-	79,79,79,79	0
57	MG	DA	1637	1/1	0.99	0.17	-	158,158,158,158	0
57	MG	AA	1606	1/1	0.98	0.04	-	102,102,102,102	0
57	MG	CA	3117	1/1	0.96	0.08	-	45,45,45,45	0
57	MG	CA	3010	1/1	0.98	0.10	-	34,34,34,34	0
57	MG	BA	3107	1/1	0.96	0.07	-	60,60,60,60	0
57	MG	BO	201	1/1	0.97	0.10	-	67,67,67,67	0
57	MG	BA	3172	1/1	0.92	0.27	-	57,57,57,57	0
57	MG	BA	3084	1/1	0.99	0.11	-	36,36,36,36	0
57	MG	CA	3007	1/1	0.97	0.28	-	204,204,204,204	0
57	MG	AA	1670	1/1	0.97	0.55	-	50,50,50,50	0
57	MG	AA	1629	1/1	0.98	0.06	-	54,54,54,54	0
57	MG	BA	3009	1/1	0.89	0.14	-	63,63,63,63	0
57	MG	CA	3103	1/1	0.86	0.04	-	116,116,116,116	0
57	MG	BA	3159	1/1	0.89	0.40	-	38,38,38,38	0
57	MG	BA	3082	1/1	0.94	0.24	-	139,139,139,139	0
57	MG	DA	1627	1/1	0.96	0.09	-	80,80,80,80	0
57	MG	CA	3027	1/1	0.98	0.04	-	47,47,47,47	0
57	MG	CA	3128	1/1	0.98	0.10	-	66,66,66,66	0
57	MG	CA	3138	1/1	0.94	0.36	-	39,39,39,39	0
57	MG	AA	1618	1/1	0.90	0.20	-	43,43,43,43	0
57	MG	AN	201	1/1	0.97	0.08	-	109,109,109,109	0
57	MG	CA	3119	1/1	0.85	0.17	-	215,215,215,215	0
57	MG	CA	3144	1/1	0.99	0.07	-	56,56,56,56	0
57	MG	DA	1640	1/1	0.96	0.30	-	66,66,66,66	0
57	MG	BA	3102	1/1	0.86	1.00	-	293,293,293,293	0
57	MG	BA	3008	1/1	0.99	0.09	-	62,62,62,62	0
57	MG	AA	1659	1/1	0.95	0.09	-	84,84,84,84	0
57	MG	DA	1653	1/1	0.97	0.20	-	51,51,51,51	0
57	MG	AA	1662	1/1	0.95	0.18	-	69,69,69,69	0
57	MG	BA	3115	1/1	0.96	0.25	-	133,133,133,133	0
57	MG	CA	3077	1/1	0.97	0.15	-	131,131,131,131	0
57	MG	CA	3056	1/1	0.97	0.23	-	155,155,155,155	0
57	MG	CA	3028	1/1	0.97	0.41	-	312,312,312,312	0
57	MG	CQ	201	1/1	0.98	0.39	-	32,32,32,32	0
57	MG	BA	3043	1/1	0.99	0.13	-	66,66,66,66	0
57	MG	DA	1615	1/1	0.98	0.20	-	213,213,213,213	0
57	MG	CA	3093	1/1	0.98	0.08	-	127,127,127,127	0
57	MG	BA	3104	1/1	0.96	0.10	-	37,37,37,37	0
57	MG	BA	3137	1/1	0.94	0.10	-	95,95,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3115	1/1	0.97	0.12	-	235,235,235,235	0
57	MG	AA	1620	1/1	0.96	0.12	-	133,133,133,133	0
57	MG	DA	1617	1/1	0.94	0.24	-	65,65,65,65	0
57	MG	AA	1651	1/1	0.89	0.27	-	60,60,60,60	0
57	MG	CA	3139	1/1	0.99	0.29	-	52,52,52,52	0
57	MG	AA	1665	1/1	0.89	0.70	-	106,106,106,106	0
57	MG	AA	1642	1/1	0.91	0.30	-	60,60,60,60	0
57	MG	AA	1661	1/1	0.93	0.14	-	53,53,53,53	0
57	MG	BA	3116	1/1	0.99	0.22	-	133,133,133,133	0
57	MG	BA	3193	1/1	0.99	0.34	-	28,28,28,28	0
57	MG	BA	3130	1/1	0.97	0.31	-	270,270,270,270	0
57	MG	BA	3012	1/1	0.90	0.07	-	100,100,100,100	0
57	MG	CA	3122	1/1	0.98	0.20	-	88,88,88,88	0
57	MG	BA	3142	1/1	0.96	0.18	-	63,63,63,63	0
57	MG	CA	3062	1/1	0.98	0.11	-	97,97,97,97	0
57	MG	BA	3143	1/1	0.98	0.35	-	25,25,25,25	0
57	MG	BA	3033	1/1	0.99	0.09	-	48,48,48,48	0
57	MG	DA	1641	1/1	0.92	0.22	-	69,69,69,69	0
57	MG	BA	3126	1/1	0.99	0.17	-	39,39,39,39	0
57	MG	CA	3043	1/1	0.94	0.07	-	57,57,57,57	0
57	MG	CA	3154	1/1	0.97	0.28	-	42,42,42,42	0
57	MG	DA	1646	1/1	0.97	0.29	-	31,31,31,31	0
57	MG	CA	3153	1/1	0.92	0.32	-	47,47,47,47	0
57	MG	DA	1623	1/1	0.97	0.07	-	64,64,64,64	0
57	MG	DD	301	1/1	0.97	0.26	-	314,314,314,314	0
57	MG	DA	1639	1/1	0.96	0.24	-	161,161,161,161	0
57	MG	BA	3071	1/1	0.97	0.11	-	43,43,43,43	0
57	MG	AA	1660	1/1	0.95	0.09	-	67,67,67,67	0
57	MG	AA	1639	1/1	0.99	0.19	-	59,59,59,59	0
57	MG	DA	1609	1/1	0.95	0.07	-	97,97,97,97	0
57	MG	BA	3093	1/1	0.93	0.11	-	72,72,72,72	0
57	MG	DA	1631	1/1	0.97	0.17	-	196,196,196,196	0
57	MG	AA	1628	1/1	0.87	0.14	-	106,106,106,106	0
57	MG	BA	3194	1/1	0.94	0.24	-	46,46,46,46	0
57	MG	AA	1637	1/1	0.89	0.23	-	138,138,138,138	0
57	MG	AA	1615	1/1	0.97	0.09	-	123,123,123,123	0
57	MG	BA	3168	1/1	0.94	0.32	-	39,39,39,39	0
57	MG	BA	3061	1/1	0.95	0.04	-	70,70,70,70	0
57	MG	BA	3191	1/1	0.99	0.13	-	43,43,43,43	0
57	MG	CA	3091	1/1	0.94	0.07	-	93,93,93,93	0
57	MG	AA	1668	1/1	0.94	0.38	-	70,70,70,70	0
57	MG	AA	1609	1/1	0.89	0.18	-	227,227,227,227	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	BA	3041	1/1	0.99	0.13	-	32,32,32,32	0
57	MG	AA	1654	1/1	0.96	0.15	-	82,82,82,82	0
57	MG	BA	3182	1/1	0.92	0.28	-	59,59,59,59	0
57	MG	AA	1647	1/1	0.99	0.20	-	44,44,44,44	0
57	MG	BA	3165	1/1	0.94	0.26	-	60,60,60,60	0
57	MG	BA	3039	1/1	0.96	0.18	-	48,48,48,48	0
57	MG	BA	3059	1/1	0.96	0.10	-	60,60,60,60	0
57	MG	BA	3170	1/1	0.94	0.10	-	51,51,51,51	0
57	MG	AA	1658	1/1	0.73	0.72	-	97,97,97,97	0
57	MG	BA	3180	1/1	0.96	0.32	-	48,48,48,48	0
57	MG	CA	3163	1/1	0.97	0.26	-	44,44,44,44	0
57	MG	BQ	201	1/1	0.86	1.37	-	108,108,108,108	0
57	MG	CA	3135	1/1	0.94	0.14	-	121,121,121,121	0
57	MG	DA	1652	1/1	0.98	0.17	-	67,67,67,67	0
57	MG	BA	3090	1/1	0.96	0.21	-	71,71,71,71	0
57	MG	CA	3162	1/1	0.91	0.30	-	71,71,71,71	0
57	MG	CA	3030	1/1	0.82	0.17	-	154,154,154,154	0
57	MG	CA	3125	1/1	0.94	0.08	-	38,38,38,38	0
57	MG	BA	3096	1/1	0.99	0.03	-	66,66,66,66	0
57	MG	BA	3131	1/1	0.99	0.08	-	39,39,39,39	0
57	MG	BA	3078	1/1	0.98	0.10	-	86,86,86,86	0
57	MG	DA	1602	1/1	0.95	0.12	-	83,83,83,83	0
57	MG	CA	3045	1/1	0.96	0.17	-	46,46,46,46	0
57	MG	AA	1638	1/1	0.97	0.08	-	58,58,58,58	0
57	MG	BA	3169	1/1	0.96	0.34	-	88,88,88,88	0
57	MG	AA	1656	1/1	0.89	0.17	-	73,73,73,73	0
57	MG	CA	3055	1/1	0.72	0.12	-	114,114,114,114	0
57	MG	AA	1634	1/1	0.97	0.08	-	87,87,87,87	0
57	MG	BA	3156	1/1	0.99	0.41	-	35,35,35,35	0
57	MG	DA	1628	1/1	0.96	0.25	-	177,177,177,177	0
57	MG	BA	3065	1/1	0.97	0.16	-	98,98,98,98	0
57	MG	CA	3061	1/1	0.89	0.26	-	102,102,102,102	0
57	MG	BA	3148	1/1	0.98	0.31	-	34,34,34,34	0
57	MG	BA	3099	1/1	0.98	0.06	-	96,96,96,96	0
57	MG	CA	3164	1/1	0.95	0.32	-	44,44,44,44	0
57	MG	BA	3123	1/1	0.95	0.06	-	115,115,115,115	0
57	MG	BA	3060	1/1	0.89	0.15	-	146,146,146,146	0
57	MG	BA	3132	1/1	0.99	0.18	-	30,30,30,30	0
57	MG	BA	3029	1/1	0.95	0.47	-	240,240,240,240	0
57	MG	AA	1641	1/1	0.97	0.11	-	124,124,124,124	0
57	MG	AA	1664	1/1	0.95	0.07	-	67,67,67,67	0
57	MG	CA	3053	1/1	0.82	0.16	-	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3127	1/1	0.98	0.06	-	47,47,47,47	0
57	MG	DD	302	1/1	0.96	0.30	-	53,53,53,53	0
57	MG	AA	1648	1/1	0.83	0.23	-	77,77,77,77	0
57	MG	CA	3089	1/1	0.97	0.16	-	215,215,215,215	0
57	MG	BA	3036	1/1	0.96	0.13	-	38,38,38,38	0
57	MG	DA	1650	1/1	0.58	0.17	-	81,81,81,81	0
57	MG	BA	3091	1/1	0.93	0.12	-	94,94,94,94	0
57	MG	BA	3119	1/1	0.90	0.24	-	262,262,262,262	0
57	MG	CA	3081	1/1	0.98	0.05	-	48,48,48,48	0
57	MG	AA	1605	1/1	0.96	0.03	-	122,122,122,122	0
57	MG	BA	3023	1/1	0.95	0.24	-	38,38,38,38	0
57	MG	CA	3021	1/1	0.98	0.05	-	57,57,57,57	0
57	MG	BA	3144	1/1	0.96	0.25	-	34,34,34,34	0
57	MG	CA	3100	1/1	0.96	0.10	-	155,155,155,155	0
57	MG	CA	3011	1/1	0.92	0.08	-	54,54,54,54	0
57	MG	BA	3037	1/1	0.97	0.17	-	52,52,52,52	0
57	MG	DA	1647	1/1	0.94	0.14	-	59,59,59,59	0
57	MG	DN	201	1/1	0.91	0.48	-	317,317,317,317	0
57	MG	BA	3192	1/1	0.99	0.18	-	46,46,46,46	0
57	MG	CA	3006	1/1	0.94	0.06	-	83,83,83,83	0
57	MG	BA	3186	1/1	0.96	0.37	-	37,37,37,37	0
57	MG	AA	1653	1/1	0.96	0.22	-	82,82,82,82	0
57	MG	BA	3175	1/1	0.96	0.18	-	44,44,44,44	0
57	MG	DA	1622	1/1	0.88	0.88	-	331,331,331,331	0
57	MG	BA	3087	1/1	0.96	0.06	-	139,139,139,139	0
57	MG	CA	3058	1/1	0.99	0.13	-	55,55,55,55	0
57	MG	AA	1646	1/1	0.41	0.45	-	99,99,99,99	0
57	MG	CA	3052	1/1	0.96	0.09	-	54,54,54,54	0
57	MG	BA	3174	1/1	0.96	0.24	-	45,45,45,45	0
57	MG	BB	3203	1/1	0.99	0.05	-	52,52,52,52	0
57	MG	CA	3111	1/1	0.98	0.44	-	164,164,164,164	0
57	MG	BA	3074	1/1	0.89	0.20	-	105,105,105,105	0
57	MG	BA	3173	1/1	0.83	0.38	-	59,59,59,59	0
57	MG	BA	3152	1/1	0.97	0.15	-	56,56,56,56	0
57	MG	CA	3092	1/1	0.88	0.20	-	134,134,134,134	0
57	MG	BA	3016	1/1	0.97	0.10	-	56,56,56,56	0
57	MG	CA	3073	1/1	0.97	0.12	-	64,64,64,64	0
57	MG	CA	3079	1/1	0.99	0.12	-	195,195,195,195	0
57	MG	DA	1604	1/1	0.89	0.38	-	166,166,166,166	0
57	MG	BA	3085	1/1	0.98	0.11	-	57,57,57,57	0
57	MG	BA	3185	1/1	0.99	0.39	-	32,32,32,32	0
57	MG	CA	3026	1/1	0.95	0.14	-	202,202,202,202	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3082	1/1	0.98	0.12	-	50,50,50,50	0
57	MG	BA	3151	1/1	0.98	0.22	-	44,44,44,44	0
57	MG	CA	3099	1/1	0.96	0.16	-	58,58,58,58	0
57	MG	DA	1624	1/1	0.97	0.11	-	30,30,30,30	0
57	MG	BA	3103	1/1	0.96	0.08	-	41,41,41,41	0
57	MG	AA	1655	1/1	0.95	0.46	-	77,77,77,77	0
57	MG	BA	3160	1/1	0.90	0.36	-	74,74,74,74	0
57	MG	BA	3190	1/1	0.89	0.23	-	68,68,68,68	0
57	MG	CA	3004	1/1	0.98	0.12	-	130,130,130,130	0
57	MG	BA	3167	1/1	0.93	0.35	-	53,53,53,53	0
57	MG	CB	3203	1/1	0.97	0.07	-	115,115,115,115	0
57	MG	DA	1651	1/1	0.96	0.20	-	51,51,51,51	0
57	MG	CA	3047	1/1	0.94	0.20	-	211,211,211,211	0
57	MG	CA	3137	1/1	0.97	0.31	-	22,22,22,22	0
57	MG	CA	3076	1/1	0.98	0.18	-	77,77,77,77	0
57	MG	BA	3150	1/1	0.98	0.71	-	38,38,38,38	0
57	MG	BA	3178	1/1	0.84	0.28	-	63,63,63,63	0
57	MG	BA	3158	1/1	0.99	0.39	-	30,30,30,30	0
57	MG	AA	1657	1/1	0.97	0.20	-	78,78,78,78	0
57	MG	CA	3001	1/1	0.97	0.09	-	54,54,54,54	0
57	MG	AA	1644	1/1	0.89	0.26	-	86,86,86,86	0
57	MG	BA	3056	1/1	0.94	0.11	-	59,59,59,59	0
57	MG	BA	3079	1/1	0.97	0.13	-	42,42,42,42	0
57	MG	CA	3123	1/1	0.93	0.25	-	177,177,177,177	0
57	MG	DA	1611	1/1	0.97	0.05	-	60,60,60,60	0
57	MG	CA	3075	1/1	0.99	0.10	-	107,107,107,107	0
57	MG	BA	3177	1/1	0.99	0.06	-	86,86,86,86	0
57	MG	CA	3114	1/1	0.99	0.13	-	169,169,169,169	0
57	MG	AA	1607	1/1	0.94	0.74	-	327,327,327,327	0
57	MG	DA	1642	1/1	0.97	0.29	-	60,60,60,60	0
57	MG	BA	3055	1/1	0.95	0.09	-	34,34,34,34	0
57	MG	CA	3086	1/1	0.91	0.23	-	83,83,83,83	0
57	MG	BA	3046	1/1	0.98	0.06	-	58,58,58,58	0
57	MG	CA	3031	1/1	0.98	0.17	-	30,30,30,30	0
57	MG	CA	3157	1/1	0.95	0.16	-	54,54,54,54	0
57	MG	AA	1627	1/1	0.91	0.21	-	111,111,111,111	0
57	MG	AA	1625	1/1	1.00	0.13	-	50,50,50,50	0
57	MG	BA	3035	1/1	0.97	0.12	-	43,43,43,43	0
57	MG	CA	3042	1/1	0.93	0.10	-	67,67,67,67	0
57	MG	CA	3040	1/1	1.00	0.13	-	98,98,98,98	0
57	MG	AA	1616	1/1	0.90	0.10	-	118,118,118,118	0
57	MG	BA	3048	1/1	0.97	0.16	-	69,69,69,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	DA	1649	1/1	0.97	0.09	-	63,63,63,63	0
57	MG	CA	3155	1/1	0.98	0.30	-	39,39,39,39	0
57	MG	BA	3073	1/1	0.99	0.09	-	141,141,141,141	0
57	MG	CA	3141	1/1	0.85	0.51	-	68,68,68,68	0
57	MG	DA	1619	1/1	0.89	0.04	-	155,155,155,155	0
57	MG	AA	1633	1/1	0.95	0.10	-	86,86,86,86	0
57	MG	BA	3088	1/1	0.97	0.09	-	147,147,147,147	0
57	MG	BA	3127	1/1	0.99	0.11	-	88,88,88,88	0
57	MG	BA	3162	1/1	0.88	0.29	-	44,44,44,44	0
57	MG	AA	1611	1/1	0.89	0.08	-	125,125,125,125	0
57	MG	CA	3078	1/1	0.89	0.05	-	113,113,113,113	0
57	MG	BA	3006	1/1	0.97	0.12	-	40,40,40,40	0
57	MG	BA	3141	1/1	0.99	0.19	-	31,31,31,31	0
57	MG	CA	3158	1/1	0.96	0.45	-	57,57,57,57	0
57	MG	BA	3050	1/1	0.99	0.10	-	27,27,27,27	0
57	MG	BA	3117	1/1	0.97	0.19	-	65,65,65,65	0
57	MG	CA	3084	1/1	0.57	0.20	-	193,193,193,193	0
57	MG	DA	1636	1/1	0.90	0.09	-	82,82,82,82	0
57	MG	CA	3126	1/1	0.92	0.16	-	146,146,146,146	0
57	MG	BA	3011	1/1	0.99	0.07	-	34,34,34,34	0
57	MG	AA	1652	1/1	0.90	0.10	-	50,50,50,50	0
57	MG	CA	3065	1/1	0.92	0.09	-	76,76,76,76	0
57	MG	CA	3071	1/1	0.92	0.51	-	218,218,218,218	0
57	MG	BA	3020	1/1	0.98	0.08	-	41,41,41,41	0
57	MG	CA	3102	1/1	0.94	0.11	-	123,123,123,123	0
57	MG	BA	3097	1/1	0.98	0.06	-	91,91,91,91	0
57	MG	BA	3153	1/1	0.97	0.35	-	34,34,34,34	0
57	MG	DA	1626	1/1	0.97	0.07	-	98,98,98,98	0
57	MG	CA	3088	1/1	0.99	0.07	-	98,98,98,98	0
57	MG	AA	1650	1/1	0.97	0.40	-	55,55,55,55	0
57	MG	CA	3038	1/1	0.93	0.10	-	71,71,71,71	0
57	MG	BL	202	1/1	0.94	0.43	-	259,259,259,259	0
57	MG	CA	3036	1/1	0.99	0.12	-	53,53,53,53	0
57	MG	BA	3040	1/1	0.97	0.09	-	66,66,66,66	0
57	MG	BA	3092	1/1	0.86	0.16	-	173,173,173,173	0
57	MG	BA	3189	1/1	0.99	0.15	-	47,47,47,47	0
57	MG	CA	3068	1/1	0.92	0.19	-	58,58,58,58	0
57	MG	CA	3069	1/1	0.97	0.17	-	63,63,63,63	0
57	MG	BA	3019	1/1	0.86	0.25	-	169,169,169,169	0
57	MG	BA	3081	1/1	0.96	0.07	-	57,57,57,57	0
57	MG	CA	3160	1/1	0.90	0.21	-	79,79,79,79	0
57	MG	CA	3095	1/1	0.79	0.77	-	270,270,270,270	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	CA	3149	1/1	0.96	0.24	-	69,69,69,69	0
57	MG	CA	3113	1/1	0.96	0.23	-	85,85,85,85	0
57	MG	BA	3110	1/1	0.95	0.44	-	89,89,89,89	0
57	MG	BA	3106	1/1	0.96	0.08	-	71,71,71,71	0
57	MG	BA	3181	1/1	0.88	0.25	-	106,106,106,106	0
57	MG	BA	3049	1/1	0.99	0.10	-	57,57,57,57	0
57	MG	CA	3057	1/1	0.99	0.08	-	46,46,46,46	0
57	MG	CA	3090	1/1	0.96	0.09	-	69,69,69,69	0
57	MG	BA	3064	1/1	0.95	0.17	-	80,80,80,80	0
57	MG	AA	1608	1/1	0.86	0.14	-	159,159,159,159	0
57	MG	CA	3020	1/1	0.99	0.14	-	37,37,37,37	0
57	MG	BA	3145	1/1	0.94	1.47	-	164,164,164,164	0
57	MG	CA	3140	1/1	0.97	0.27	-	29,29,29,29	0
57	MG	BA	3129	1/1	0.97	0.19	-	78,78,78,78	0
57	MG	BA	3164	1/1	0.94	0.31	-	54,54,54,54	0
57	MG	BB	3204	1/1	0.99	0.20	-	24,24,24,24	0
57	MG	AA	1613	1/1	0.98	0.10	-	97,97,97,97	0
57	MG	CA	3003	1/1	0.97	0.07	-	95,95,95,95	0
57	MG	BB	3202	1/1	0.98	0.08	-	34,34,34,34	0
57	MG	CA	3034	1/1	0.95	0.29	-	157,157,157,157	0
57	MG	CA	3044	1/1	0.98	0.05	-	53,53,53,53	0
57	MG	BA	3086	1/1	0.96	0.12	-	74,74,74,74	0
57	MG	BA	3098	1/1	0.98	0.45	-	157,157,157,157	0
57	MG	BA	3100	1/1	0.98	0.15	-	66,66,66,66	0
57	MG	BA	3139	1/1	0.98	0.12	-	47,47,47,47	0
57	MG	AA	1643	1/1	0.91	0.70	-	162,162,162,162	0
57	MG	BA	3080	1/1	0.95	0.19	-	75,75,75,75	0
57	MG	CA	3148	1/1	0.97	0.18	-	37,37,37,37	0
57	MG	BA	3063	1/1	0.98	0.24	-	93,93,93,93	0
57	MG	BA	3030	1/1	0.98	0.08	-	43,43,43,43	0
57	MG	CA	3029	1/1	0.98	0.37	-	168,168,168,168	0
57	MG	CA	3070	1/1	0.97	0.14	-	251,251,251,251	0
57	MG	CA	3165	1/1	0.92	0.38	-	83,83,83,83	0
57	MG	BA	3076	1/1	0.98	0.10	-	32,32,32,32	0
57	MG	BA	3171	1/1	1.00	0.19	-	47,47,47,47	0
57	MG	BA	3121	1/1	0.98	0.10	-	34,34,34,34	0
57	MG	DA	1634	1/1	0.79	0.06	-	128,128,128,128	0
57	MG	DA	1608	1/1	0.77	0.16	-	218,218,218,218	0
57	MG	CA	3085	1/1	0.93	0.33	-	260,260,260,260	0
57	MG	CA	3046	1/1	0.98	0.27	-	130,130,130,130	0
57	MG	AA	1666	1/1	0.96	0.23	-	44,44,44,44	0
57	MG	DA	1635	1/1	0.97	0.24	-	126,126,126,126	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
57	MG	AA	1671	1/1	0.91	0.25	-	47,47,47,47	0
57	MG	DA	1648	1/1	0.95	0.50	-	41,41,41,41	0
57	MG	CA	3014	1/1	0.93	0.12	-	82,82,82,82	0
57	MG	BA	3146	1/1	0.99	0.31	-	48,48,48,48	0
57	MG	CA	3146	1/1	0.94	0.14	-	52,52,52,52	0
57	MG	AA	1649	1/1	0.97	0.26	-	51,51,51,51	0
57	MG	CA	3022	1/1	0.98	0.09	-	59,59,59,59	0
57	MG	BA	3052	1/1	0.96	0.16	-	50,50,50,50	0
57	MG	CA	3147	1/1	0.90	0.38	-	77,77,77,77	0
57	MG	CA	3009	1/1	0.96	0.15	-	88,88,88,88	0
57	MG	DA	1645	1/1	0.91	0.23	-	57,57,57,57	0
57	MG	BA	3018	1/1	0.95	0.13	-	59,59,59,59	0
57	MG	CA	3035	1/1	0.98	0.11	-	35,35,35,35	0
57	MG	DA	1606	1/1	0.78	0.08	-	144,144,144,144	0
57	MG	BA	3058	1/1	0.99	0.12	-	25,25,25,25	0
57	MG	AA	1632	1/1	0.95	0.18	-	149,149,149,149	0
57	MG	BA	3122	1/1	0.96	0.06	-	60,60,60,60	0
57	MG	DA	1643	1/1	0.97	0.16	-	73,73,73,73	0
57	MG	CA	3016	1/1	0.98	0.06	-	40,40,40,40	0
57	MG	AA	1663	1/1	0.92	0.32	-	52,52,52,52	0
57	MG	BA	3015	1/1	0.96	0.10	-	58,58,58,58	0
57	MG	BA	3024	1/1	0.97	0.08	-	68,68,68,68	0
57	MG	CA	3159	1/1	0.97	0.17	-	52,52,52,52	0
57	MG	BA	3051	1/1	0.98	0.10	-	70,70,70,70	0
57	MG	CA	3087	1/1	0.91	0.20	-	75,75,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.