



wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 08:27 am BST

PDB ID : 4XEJ
Title : IRES bound to bacterial Ribosome
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Deposited on : 2014-12-23
Resolution : 3.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

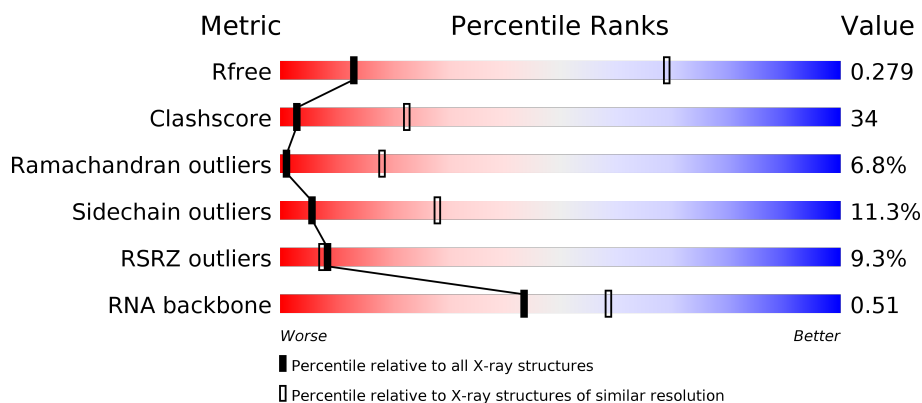
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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}	130704	1212 (4.00-3.60)
Clashscore	141614	1288 (4.00-3.60)
Ramachandran outliers	138981	1243 (4.00-3.60)
Sidechain outliers	138945	1237 (4.00-3.60)
RSRZ outliers	127900	1121 (4.00-3.60)
RNA backbone	3102	1036 (4.60-3.00)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AL02	271	<div> <div>3%</div> <div>82%</div> <div>17%</div> <div>•</div> </div>
1	BL02	271	<div> <div>77%</div> <div>21%</div> <div>•</div> </div>
2	AL03	204	<div> <div>3%</div> <div>80%</div> <div>19%</div> <div>•</div> </div>
2	BL03	204	<div> <div>3%</div> <div>77%</div> <div>21%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
3	AL04	202	
3	BL04	202	
4	AL05	181	
4	BL05	181	
5	AL06	159	
5	BL06	159	
6	AL09	145	
6	BL09	145	
7	AL11	147	
7	BL11	147	
8	AL13	137	
8	BL13	137	
9	AL14	122	
9	BL14	122	
10	AL15	146	
10	BL15	146	
11	AL16	134	
11	BL16	134	
12	AL17	117	
12	BL17	117	
13	AL18	98	
13	BL18	98	
14	AL19	137	
14	BL19	137	
15	AL20	117	

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Mol	Chain	Length	Quality of chain
15	BL20	117	
16	AL21	101	
16	BL21	101	
17	AL22	112	
17	BL22	112	
18	AL23	92	
18	BL23	92	
19	AL24	100	
19	BL24	100	
20	AL25	187	
20	BL25	187	
21	AL27	76	
21	BL27	76	
22	AL28	88	
22	BL28	88	
23	AL29	62	
23	BL29	62	
24	AL30	59	
24	BL30	59	
25	AL32	52	
25	BL32	52	
26	AL33	44	
26	BL33	44	
27	AL34	48	
27	BL34	48	

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Mol	Chain	Length	Quality of chain
28	AL35	63	
28	BL35	63	
29	AS02	234	
29	BS02	234	
30	AS03	206	
30	BS03	206	
31	AS04	208	
31	BS04	208	
32	AS05	151	
32	BS05	151	
33	AS06	101	
33	BS06	101	
34	AS07	155	
34	BS07	155	
35	AS08	138	
35	BS08	138	
36	AS09	127	
36	BS09	127	
37	AS10	98	
37	BS10	98	
38	AS11	114	
38	BS11	114	
39	AS12	122	
39	BS12	122	
40	AS13	117	

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Mol	Chain	Length	Quality of chain
40	BS13	117	<div> <div>26%</div> <div>92%</div> <div>8%</div> </div>
41	AS14	60	<div> <div>18%</div> <div>93%</div> <div>5%</div> </div>
41	BS14	60	<div> <div>15%</div> <div>92%</div> <div>8%</div> </div>
42	AS15	88	<div> <div>6%</div> <div>89%</div> <div>11%</div> </div>
42	BS15	88	<div> <div>6%</div> <div>86%</div> <div>14%</div> </div>
43	AS16	83	<div> <div>14%</div> <div>83%</div> <div>17%</div> </div>
43	BS16	83	<div> <div>33%</div> <div>93%</div> <div>7%</div> </div>
44	AS17	99	<div> <div>3%</div> <div>86%</div> <div>14%</div> </div>
44	BS17	99	<div> <div>3%</div> <div>89%</div> <div>10%</div> </div>
45	AS18	70	<div> <div>13%</div> <div>90%</div> <div>10%</div> </div>
45	BS18	70	<div> <div>%</div> <div>84%</div> <div>16%</div> </div>
46	AS19	78	<div> <div>49%</div> <div>87%</div> <div>12%</div> </div>
46	BS19	78	<div> <div>36%</div> <div>85%</div> <div>14%</div> </div>
47	AS20	99	<div> <div>9%</div> <div>86%</div> <div>14%</div> </div>
47	BS20	99	<div> <div>22%</div> <div>92%</div> <div>8%</div> </div>
48	ATHX	24	<div> <div>4%</div> <div>96%</div> <div>.</div> </div>
48	BTHX	24	<div> <div>33%</div> <div>92%</div> <div>8%</div> </div>
49	AL31	30	<div> <div>43%</div> <div>90%</div> <div>10%</div> </div>
49	BL31	30	<div> <div>37%</div> <div>87%</div> <div>10%</div> </div>
50	A16S	1506	<div> <div>3%</div> <div>83%</div> <div>17%</div> </div>
50	B16S	1506	<div> <div>4%</div> <div>83%</div> <div>17%</div> </div>
51	A23S	2879	<div> <div>4%</div> <div>81%</div> <div>19%</div> </div>
51	B23S	2879	<div> <div>4%</div> <div>80%</div> <div>20%</div> </div>
52	A5S	119	<div> <div>%</div> <div>27%</div> <div>61%</div> <div>13%</div> </div>
52	B5S	119	<div> <div>%</div> <div>25%</div> <div>66%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
53	AIRE	196	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>5%14%84%</div></div>
53	BIRE	196	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>%14%84%</div></div>

2 Entry composition [i](#)

There are 54 unique types of molecules in this entry. The entry contains 287293 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AL02	271	Total	C	N	O	S	0	0	0
			2105	1329	416	357	3			
1	BL02	271	Total	C	N	O	S	0	0	0
			2105	1329	416	357	3			

- Molecule 2 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AL03	204	Total	C	N	O	S	0	0	0
			1564	988	299	271	6			
2	BL03	204	Total	C	N	O	S	0	0	0
			1564	988	299	271	6			

- Molecule 3 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AL04	202	Total	C	N	O	S	0	0	0
			1587	1011	297	276	3			
3	BL04	202	Total	C	N	O	S	0	0	0
			1587	1011	297	276	3			

- Molecule 4 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AL05	181	Total	C	N	O	S	0	0	0
			1475	943	268	260	4			
4	BL05	181	Total	C	N	O	S	0	0	0
			1475	943	268	260	4			

- Molecule 5 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AL06	159	Total	C	N	O	S	0	0	0
			1223	773	228	221	1			
5	BL06	159	Total	C	N	O	S	0	0	0
			1223	773	228	221	1			

- Molecule 6 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AL09	145	Total	C	N	O	S	0	0	0
			1133	724	200	208	1			
6	BL09	145	Total	C	N	O	S	0	0	0
			1133	724	200	208	1			

- Molecule 7 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AL11	147	Total	C	N	O	S	0	0	0
			1088	692	191	199	6			
7	BL11	147	Total	C	N	O	S	0	0	0
			1088	692	191	199	6			

- Molecule 8 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AL13	137	Total	C	N	O	S	0	0	0
			1097	707	205	182	3			
8	BL13	137	Total	C	N	O	S	0	0	0
			1097	707	205	182	3			

- Molecule 9 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	AL14	122	Total	C	N	O	S	0	0	0
			932	587	171	170	4			
9	BL14	122	Total	C	N	O	S	0	0	0
			932	587	171	170	4			

- Molecule 10 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AL15	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	BL15	146	Total	C	N	O	S	0	0	0
			1114	692	227	193	2			

- Molecule 11 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AL16	134	Total	C	N	O	S	0	0	0
			1065	680	201	179	5			
11	BL16	134	Total	C	N	O	S	0	0	0
			1065	680	201	179	5			

- Molecule 12 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AL17	117	Total	C	N	O		0	0	0
			960	599	202	159				
12	BL17	117	Total	C	N	O		0	0	0
			960	599	202	159				

- Molecule 13 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AL18	98	Total	C	N	O		0	0	0
			771	486	154	131				
13	BL18	98	Total	C	N	O		0	0	0
			771	486	154	131				

- Molecule 14 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AL19	137	Total	C	N	O	S	0	0	0
			1144	713	234	196	1			
14	BL19	137	Total	C	N	O	S	0	0	0
			1144	713	234	196	1			

- Molecule 15 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AL20	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			
15	BL20	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			

- Molecule 16 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AL21	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			
16	BL21	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			

- Molecule 17 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AL22	112	Total	C	N	O	S	0	0	0
			891	560	175	154	2			
17	BL22	112	Total	C	N	O	S	0	0	0
			891	560	175	154	2			

- Molecule 18 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	AL23	92	Total	C	N	O	0	0	0
			726	471	131	124			
18	BL23	92	Total	C	N	O	0	0	0
			726	471	131	124			

- Molecule 19 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AL24	100	Total	C	N	O	S	0	0	0
			776	500	148	124	4			
19	BL24	100	Total	C	N	O	S	0	0	0
			776	500	148	124	4			

- Molecule 20 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AL25	187	Total	C	N	O	S	0	0	0
			1483	945	264	272	2			
20	BL25	187	Total	C	N	O	S	0	0	0
			1483	945	264	272	2			

- Molecule 21 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	AL27	76	Total	C	N	O	S	0	0	0
			605	376	126	102	1			
21	BL27	76	Total	C	N	O	S	0	0	0
			605	376	126	102	1			

- Molecule 22 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	AL28	88	Total	C	N	O	S	0	0	0
			695	435	141	119				
22	BL28	88	Total	C	N	O	S	0	0	0
			695	435	141	119				

- Molecule 23 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	AL29	62	Total	C	N	O	S	0	0	0
			521	325	102	92	2			
23	BL29	62	Total	C	N	O	S	0	0	0
			521	325	102	92	2			

- Molecule 24 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	AL30	59	Total	C	N	O	S	0	0	0
			468	298	90	79	1			
24	BL30	59	Total	C	N	O	S	0	0	0
			468	298	90	79	1			

- Molecule 25 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	AL32	52	Total	C	N	O	S	0	0	0
			405	255	79	66	5			
25	BL32	52	Total	C	N	O	S	0	0	0
			405	255	79	66	5			

- Molecule 26 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	AL33	44	Total	C	N	O	S	0	0	0
			381	235	77	65	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	BL33	44	Total	C	N	O	S	0	0	0
			381	235	77	65	4			

- Molecule 27 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	AL34	48	Total	C	N	O	S	0	0	0
			419	257	104	56	2			
27	BL34	48	Total	C	N	O	S	0	0	0
			419	257	104	56	2			

- Molecule 28 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	AL35	63	Total	C	N	O	S	0	0	0
			508	326	101	79	2			
28	BL35	63	Total	C	N	O	S	0	0	0
			508	326	101	79	2			

- Molecule 29 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	AS02	234	Total	C	N	O	S	0	0	0
			1901	1213	341	342	5			
29	BS02	234	Total	C	N	O	S	0	0	0
			1901	1213	341	342	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	AS03	206	Total	C	N	O	S	0	0	0
			1613	1016	314	282	1			
30	BS03	206	Total	C	N	O	S	0	0	0
			1613	1016	314	282	1			

- Molecule 31 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	AS04	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			
31	BS04	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

- Molecule 32 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	AS05	151	Total	C	N	O	S	0	0	0
			1156	729	218	205	4			
32	BS05	151	Total	C	N	O	S	0	0	0
			1156	729	218	205	4			

- Molecule 33 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	AS06	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			
33	BS06	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

- Molecule 34 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	AS07	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			
34	BS07	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

- Molecule 35 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	AS08	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			
35	BS08	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

- Molecule 36 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
36	AS09	127	Total	C	N	O	0	0	0
			1011	639	198	174			
36	BS09	127	Total	C	N	O	0	0	0
			1011	639	198	174			

- Molecule 37 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	AS10	98	Total	C	N	O	S	0	0	0
			795	499	156	139	1			
37	BS10	98	Total	C	N	O	S	0	0	0
			795	499	156	139	1			

- Molecule 38 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	AS11	114	Total	C	N	O	S	0	0	0
			843	522	159	159	3			
38	BS11	114	Total	C	N	O	S	0	0	0
			843	522	159	159	3			

- Molecule 39 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	AS12	122	Total	C	N	O	S	0	0	0
			957	603	193	160	1			
39	BS12	122	Total	C	N	O	S	0	0	0
			957	603	193	160	1			

- Molecule 40 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	AS13	117	Total	C	N	O	S	0	0	0
			934	577	192	163	2			
40	BS13	117	Total	C	N	O	S	0	0	0
			934	577	192	163	2			

- Molecule 41 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	AS14	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			
41	BS14	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

- Molecule 42 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	AS15	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BS15	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

- Molecule 43 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	AS16	83	Total	C	N	O	S	0	0	0
			701	443	139	118	1			
43	BS16	83	Total	C	N	O	S	0	0	0
			701	443	139	118	1			

- Molecule 44 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	AS17	99	Total	C	N	O	S	0	0	0
			824	528	152	142	2			
44	BS17	99	Total	C	N	O	S	0	0	0
			824	528	152	142	2			

- Molecule 45 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
45	AS18	70	Total	C	N	O	0	0	0
			574	367	112	95			
45	BS18	70	Total	C	N	O	0	0	0
			574	367	112	95			

- Molecule 46 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	AS19	78	Total	C	N	O	S	0	0	0
			630	403	114	111	2			
46	BS19	78	Total	C	N	O	S	0	0	0
			630	403	114	111	2			

- Molecule 47 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	AS20	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			
47	BS20	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			

- Molecule 48 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
48	ATHX	24	Total	C	N	O	0	0	0
			209	128	50	31			
48	BTHX	24	Total	C	N	O	0	0	0
			209	128	50	31			

- Molecule 49 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	AL31	30	Total	C	N	O	S	0	0	0
			226	142	36	44	4			
49	BL31	30	Total	C	N	O	S	0	0	0
			226	142	36	44	4			

- Molecule 50 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	A16S	1504	Total	C	N	O	P	0	0	0
			32332	14391	5994	10444	1503			
50	B16S	1504	Total	C	N	O	P	0	0	0
			32331	14391	5994	10443	1503			

- Molecule 51 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	A23S	2876	Total	C	N	O	P	0	0	0
			61929	27563	11567	19924	2875			
51	B23S	2876	Total	C	N	O	P	0	0	0
			61931	27565	11569	19922	2875			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A23S	1141A	U	C	conflict	GB 46197919
A23S	2825	U	G	conflict	GB 46197919
B23S	1141A	U	C	conflict	GB 46197919
B23S	2825	U	G	conflict	GB 46197919

- Molecule 52 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	A5S	119	Total 2551	C 1136	N 471	O 826	P 118	0	0	0
52	B5S	119	Total 2551	C 1136	N 471	O 826	P 118	0	0	0

- Molecule 53 is a RNA chain called IRES RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	AIRE	32	Total 672	C 302	N 114	O 225	P 31	0	0	0
53	BIRE	32	Total 672	C 302	N 114	O 225	P 31	0	0	0

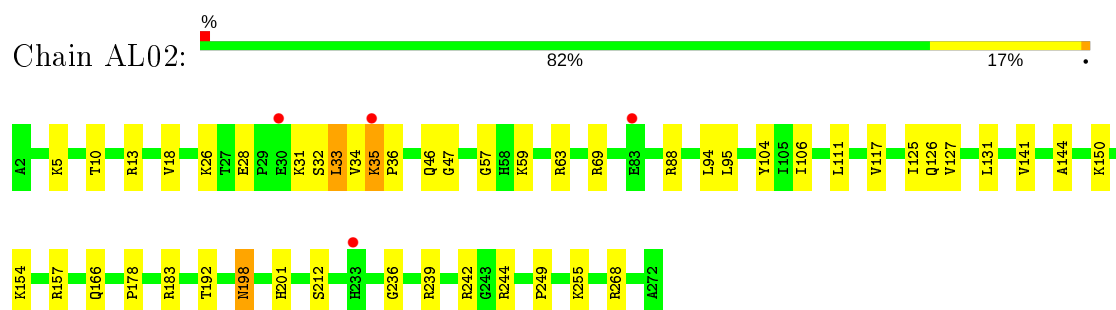
- Molecule 54 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
54	AS14	1	Total 1	Zn 1	0	0
54	AS04	1	Total 1	Zn 1	0	0
54	BS04	1	Total 1	Zn 1	0	0
54	BS14	1	Total 1	Zn 1	0	0

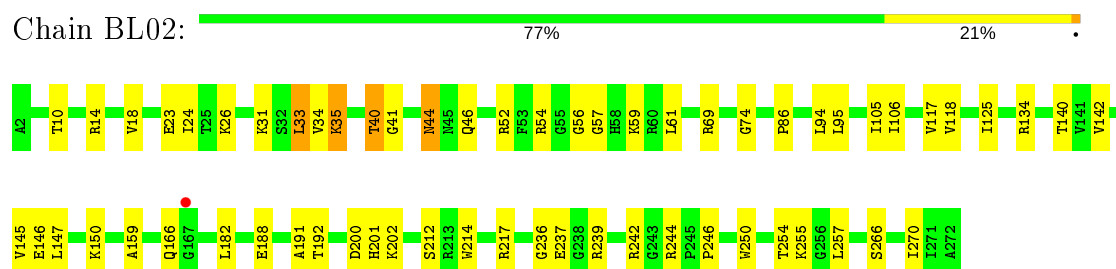
3 Residue-property plots [i](#)

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

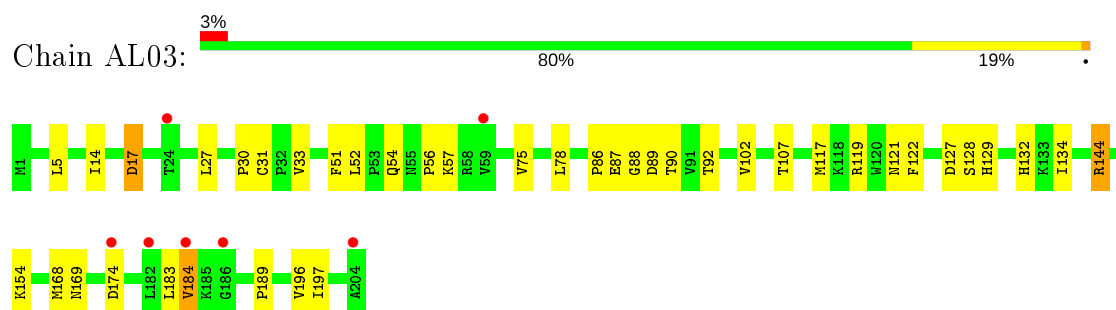
- Molecule 1: 50S ribosomal protein L2



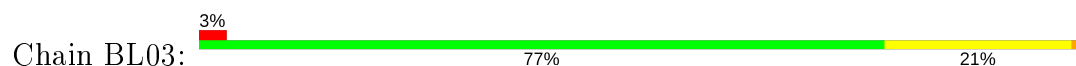
- Molecule 1: 50S ribosomal protein L2

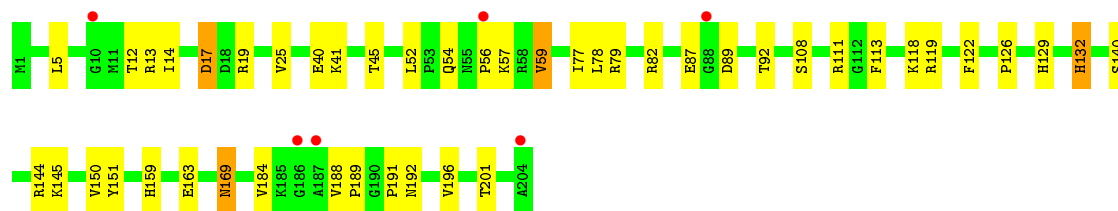


- Molecule 2: 50S ribosomal protein L3

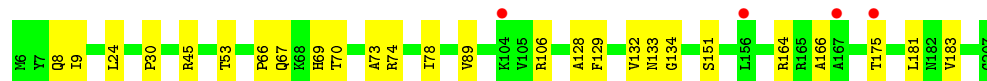


- Molecule 2: 50S ribosomal protein L3

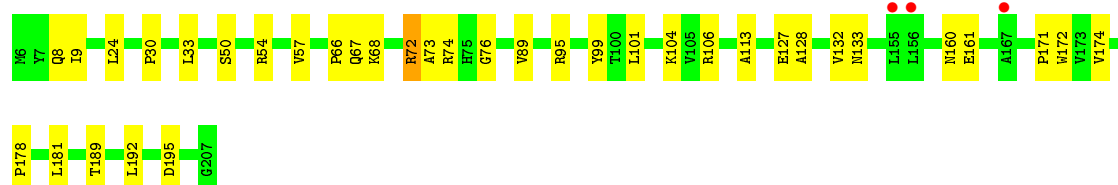
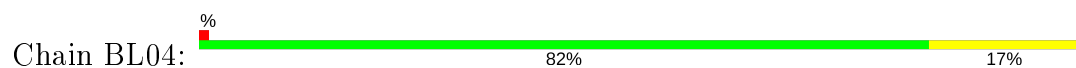




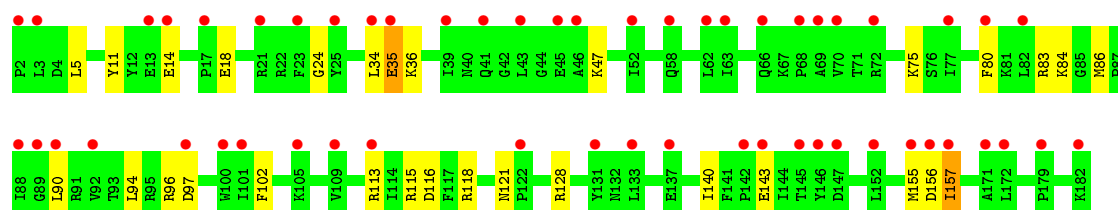
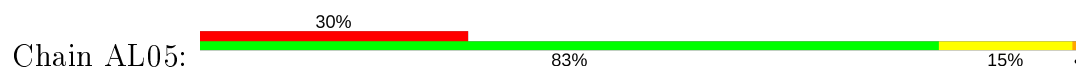
- Molecule 3: 50S ribosomal protein L4



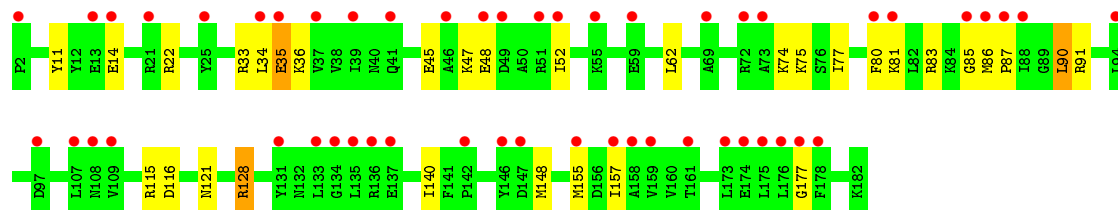
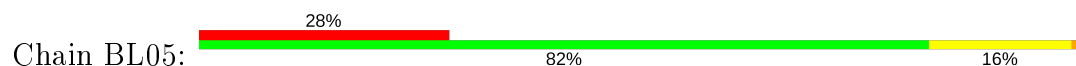
- Molecule 3: 50S ribosomal protein L4



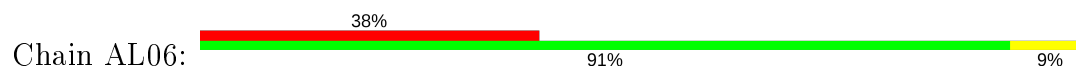
- Molecule 4: 50S ribosomal protein L5

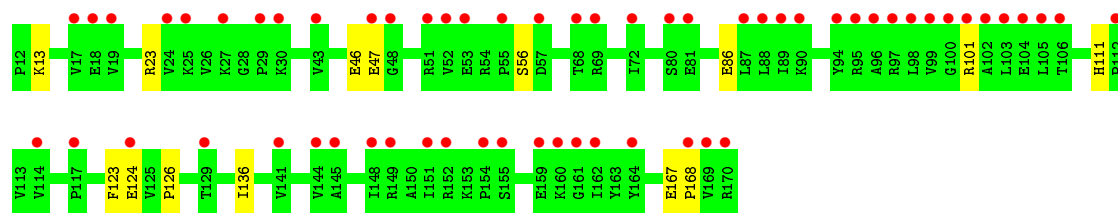


- Molecule 4: 50S ribosomal protein L5

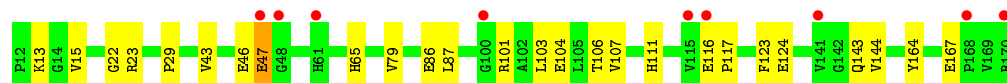
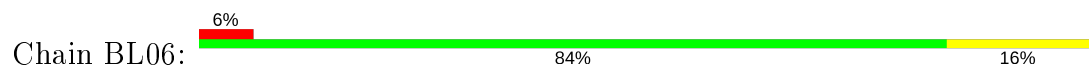


- Molecule 5: 50S ribosomal protein L6

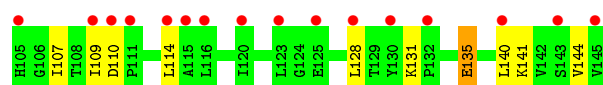
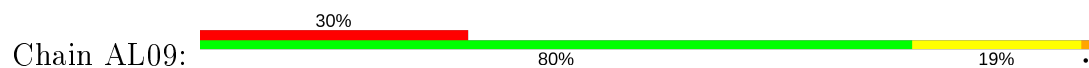




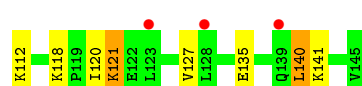
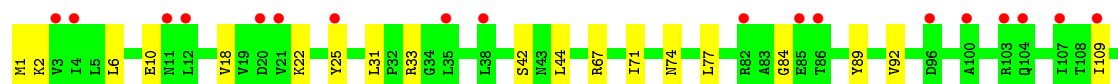
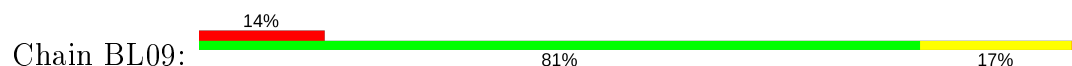
• Molecule 5: 50S ribosomal protein L6



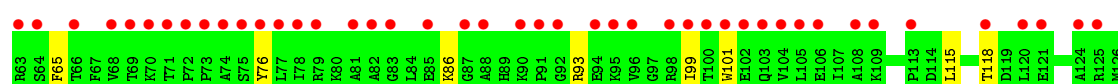
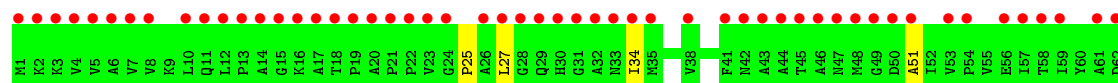
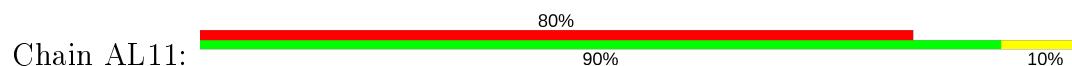
• Molecule 6: 50S ribosomal protein L9



• Molecule 6: 50S ribosomal protein L9

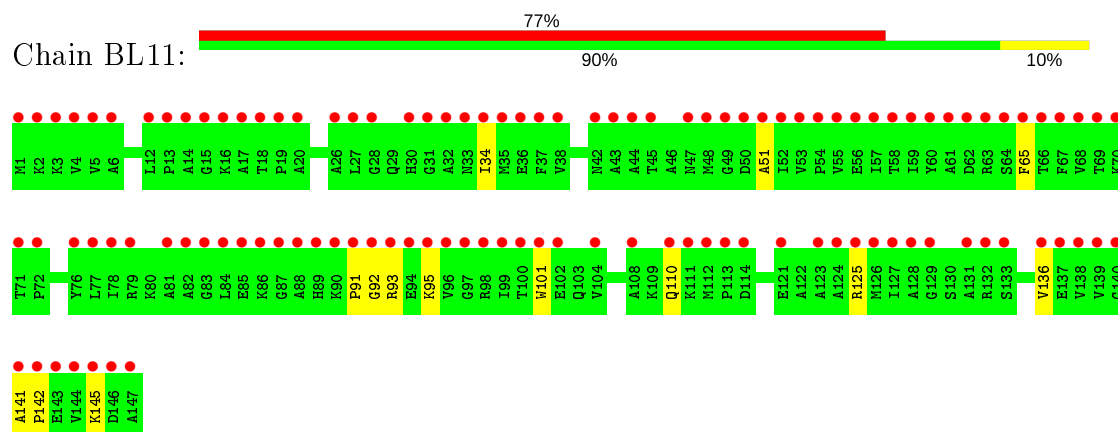


• Molecule 7: 50S ribosomal protein L11



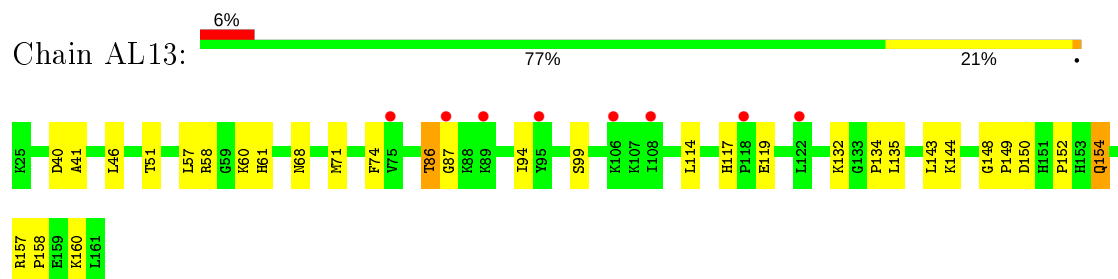
- Molecule 7: 50S ribosomal protein L11

Chain BL11:



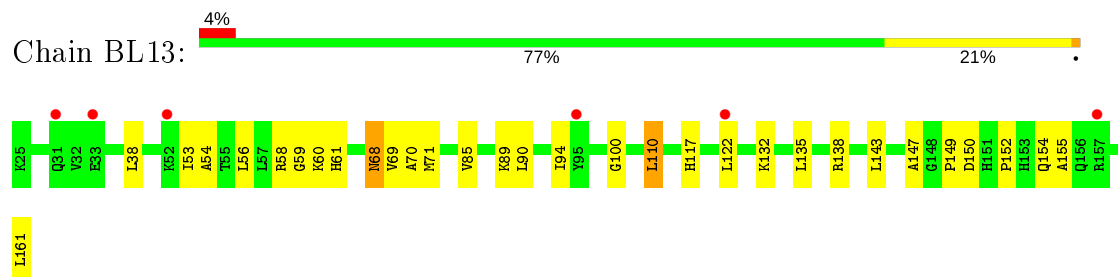
- Molecule 8: 50S ribosomal protein L13

Chain AL13:



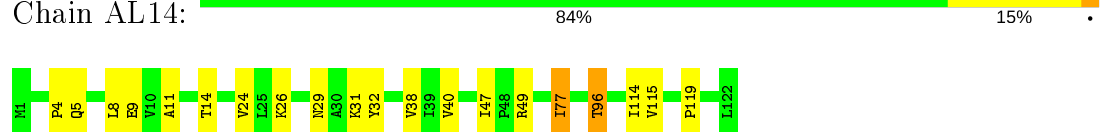
- Molecule 8: 50S ribosomal protein L13

Chain BL13:



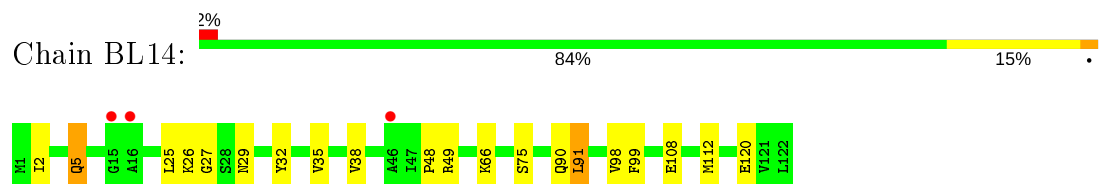
- Molecule 9: 50S ribosomal protein L14

Chain AL14:

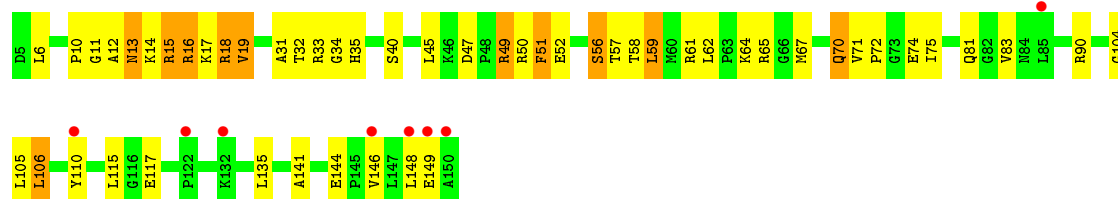


- Molecule 9: 50S ribosomal protein L14

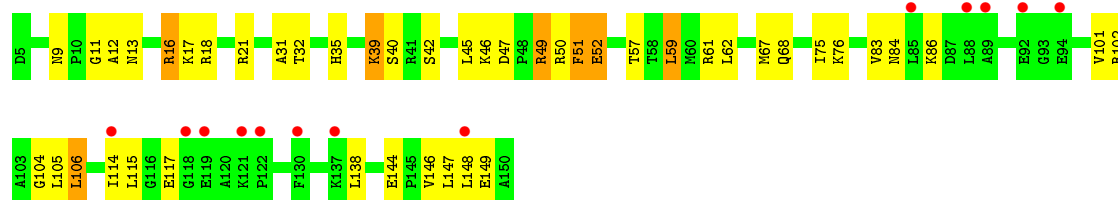
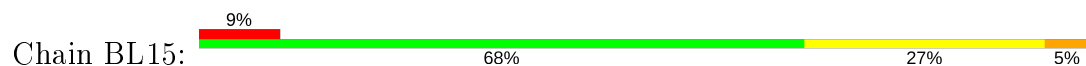
Chain BL14:



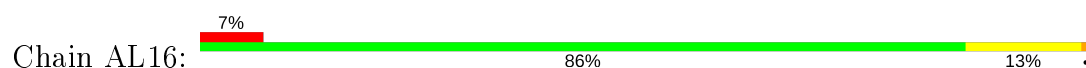
- Molecule 10: 50S ribosomal protein L15



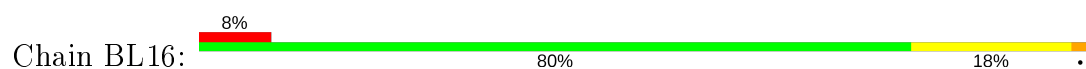
- Molecule 10: 50S ribosomal protein L15



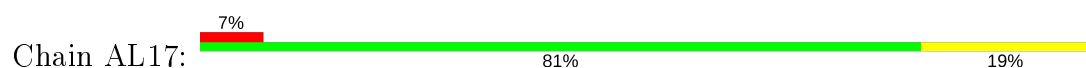
- Molecule 11: 50S ribosomal protein L16



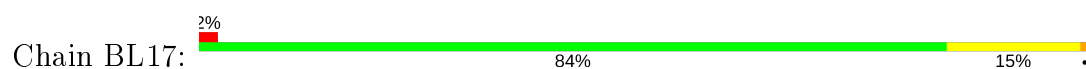
- Molecule 11: 50S ribosomal protein L16



- Molecule 12: 50S ribosomal protein L17

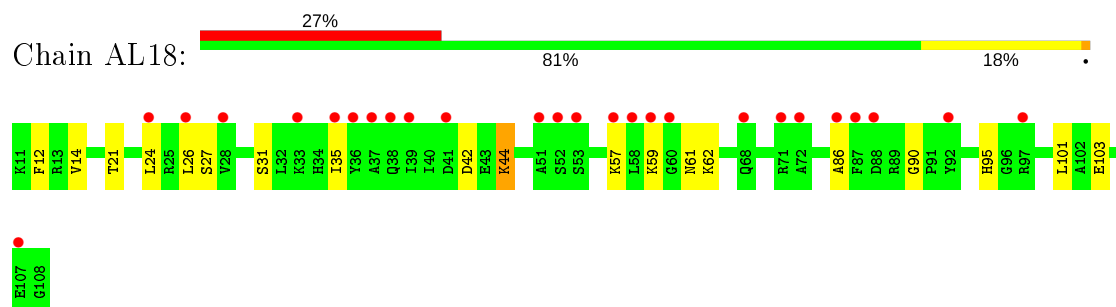


- Molecule 12: 50S ribosomal protein L17



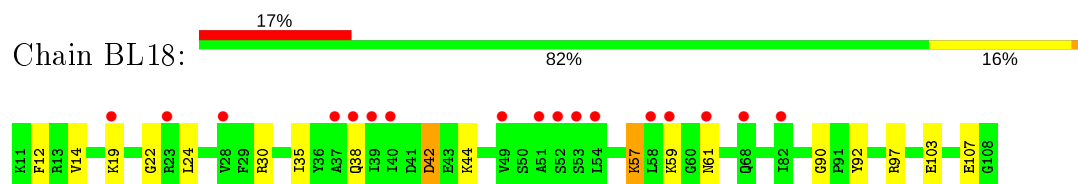
- Molecule 13: 50S ribosomal protein L18

Chain AL18:



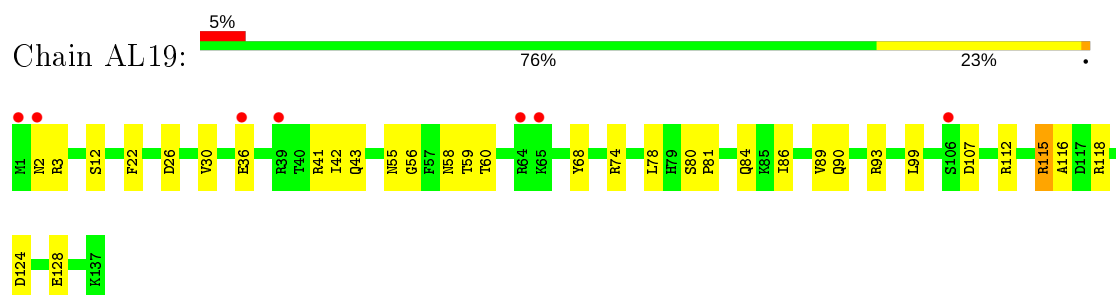
- Molecule 13: 50S ribosomal protein L18

Chain BL18:



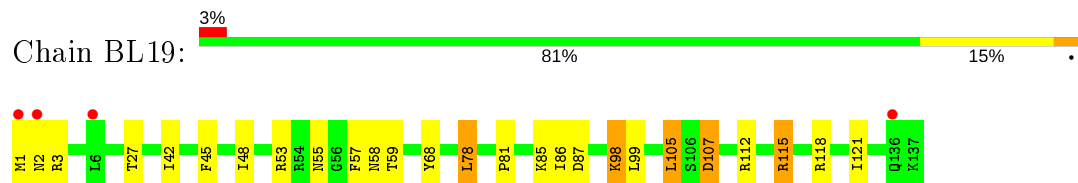
- Molecule 14: 50S ribosomal protein L19

Chain AL19:



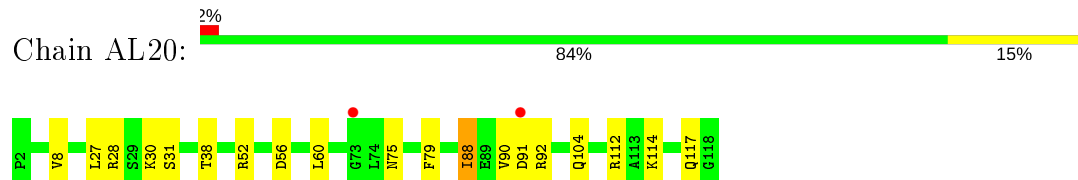
- Molecule 14: 50S ribosomal protein L19

Chain BL19:



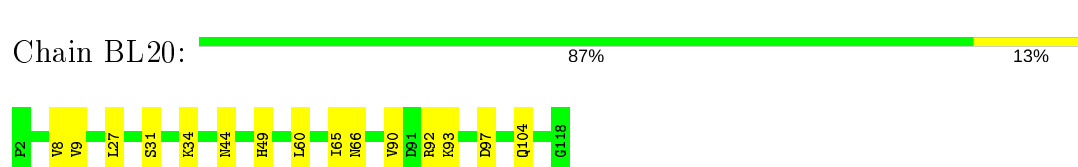
- Molecule 15: 50S ribosomal protein L20

Chain AL20:

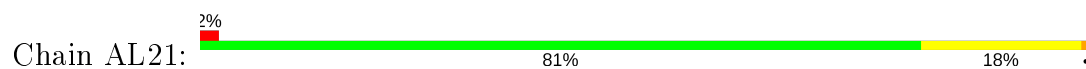


- Molecule 15: 50S ribosomal protein L20

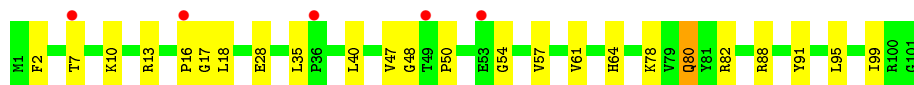
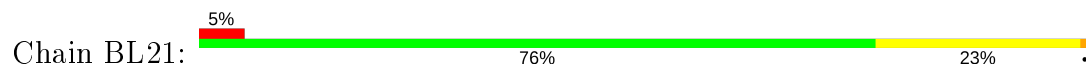
Chain BL20:



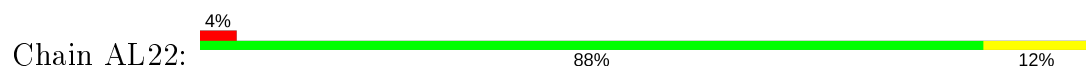
- Molecule 16: 50S ribosomal protein L21



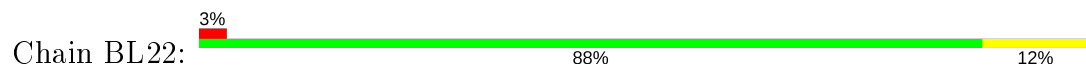
- Molecule 16: 50S ribosomal protein L21



- Molecule 17: 50S ribosomal protein L22



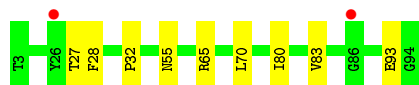
- Molecule 17: 50S ribosomal protein L22



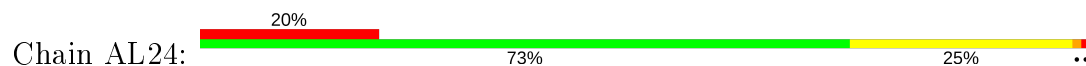
- Molecule 18: 50S ribosomal protein L23

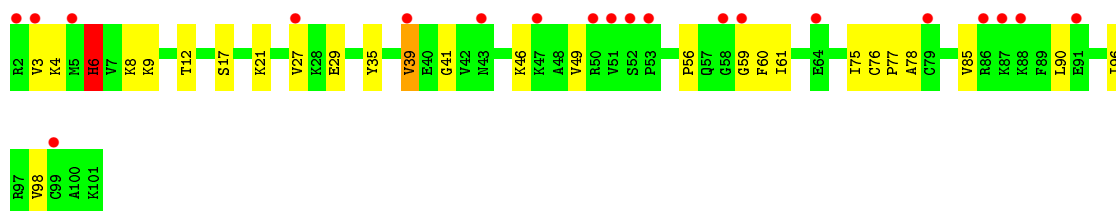


- Molecule 18: 50S ribosomal protein L23

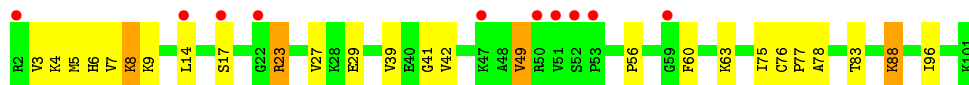
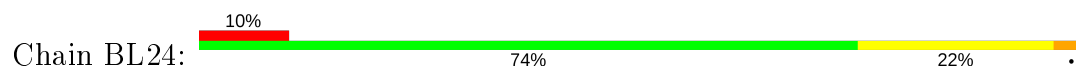


- Molecule 19: 50S ribosomal protein L24

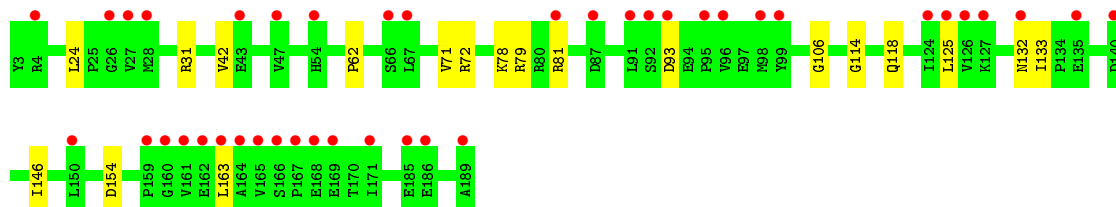




- Molecule 19: 50S ribosomal protein L24



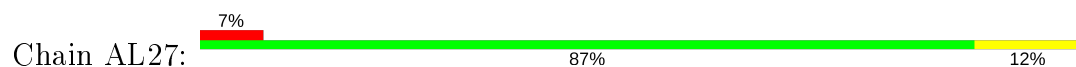
- Molecule 20: 50S ribosomal protein L25



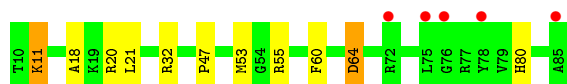
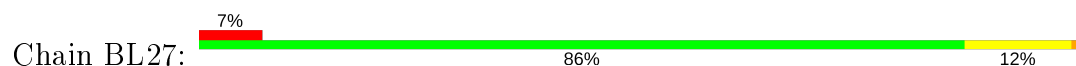
- Molecule 20: 50S ribosomal protein L25



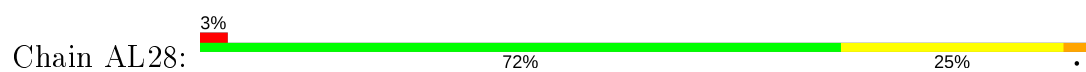
- Molecule 21: 50S ribosomal protein L27



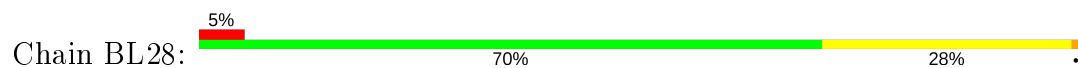
- Molecule 21: 50S ribosomal protein L27



- Molecule 22: 50S ribosomal protein L28



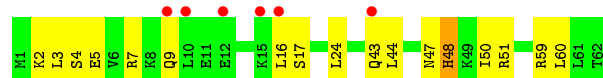
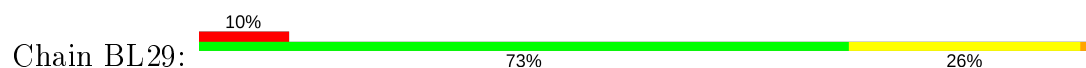
- Molecule 22: 50S ribosomal protein L28



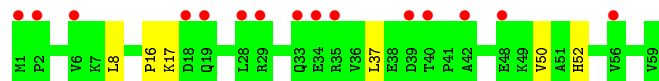
- Molecule 23: 50S ribosomal protein L29



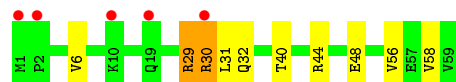
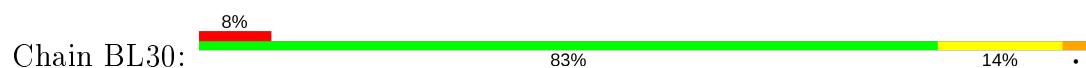
- Molecule 23: 50S ribosomal protein L29



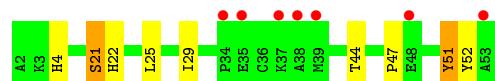
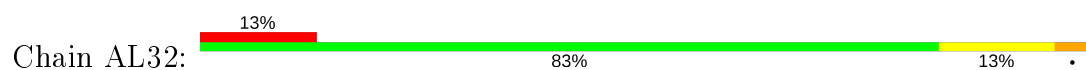
- Molecule 24: 50S ribosomal protein L30



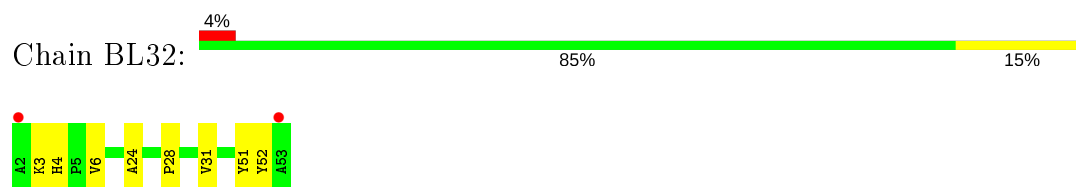
- Molecule 24: 50S ribosomal protein L30



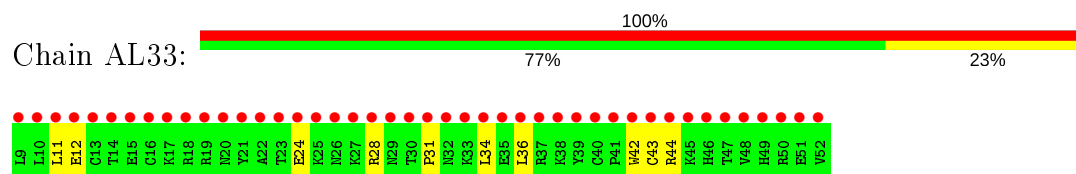
- Molecule 25: 50S ribosomal protein L32



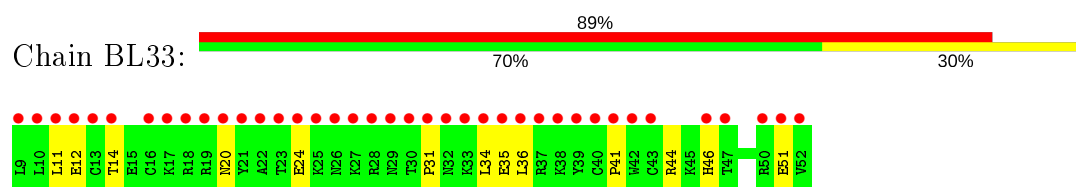
- Molecule 25: 50S ribosomal protein L32



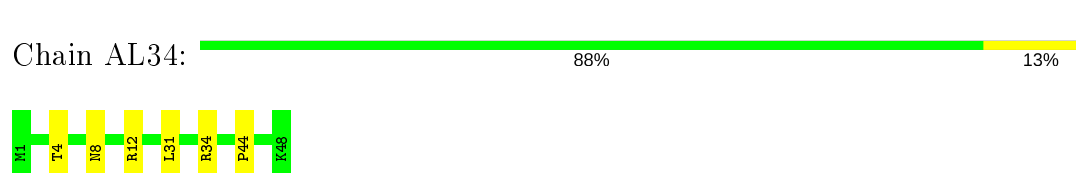
- Molecule 26: 50S ribosomal protein L33



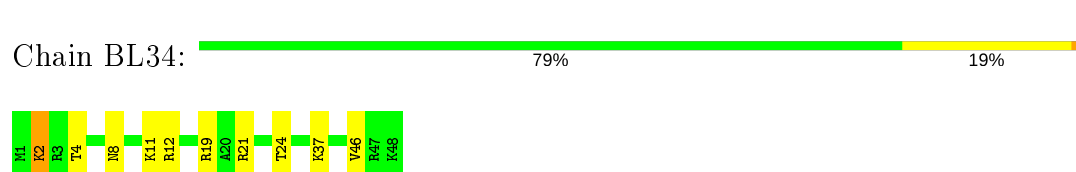
- Molecule 26: 50S ribosomal protein L33



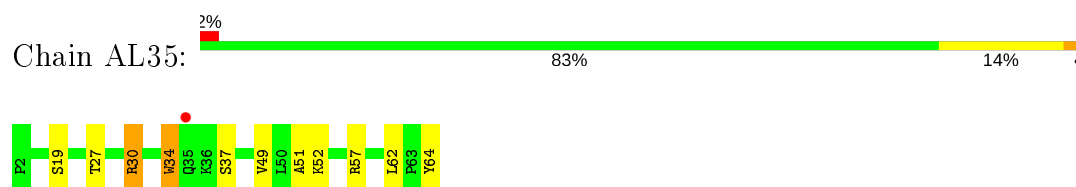
- Molecule 27: 50S ribosomal protein L34



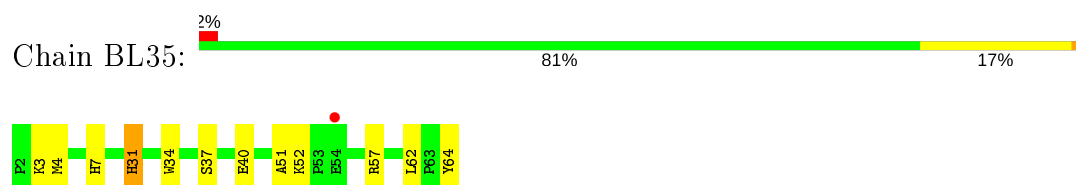
- Molecule 27: 50S ribosomal protein L34



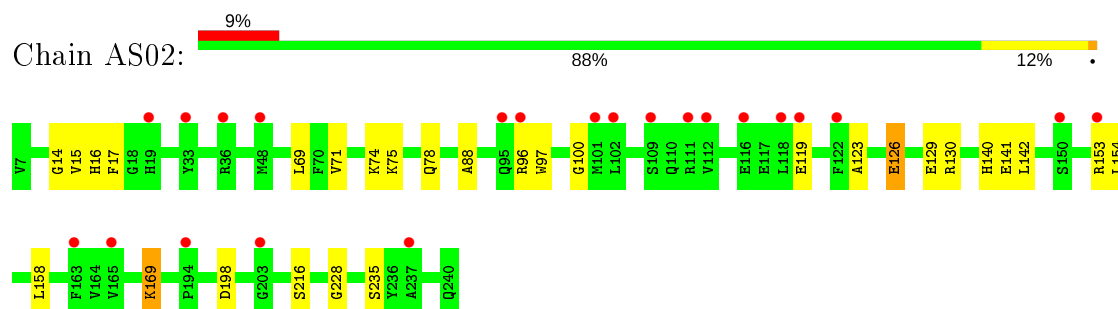
- Molecule 28: 50S ribosomal protein L35



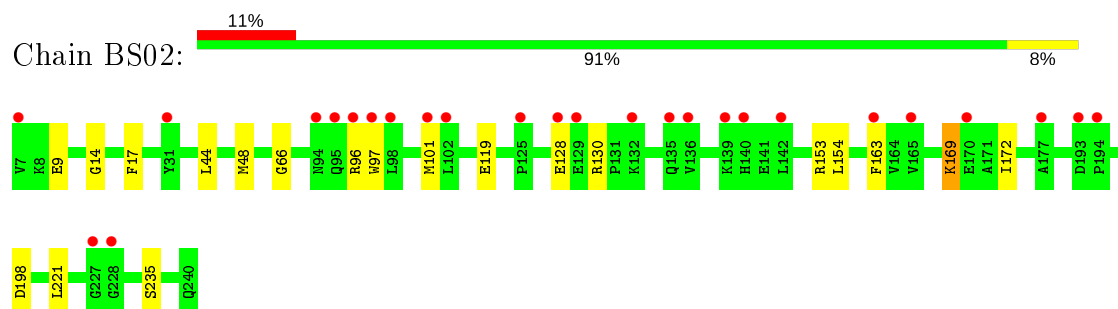
- Molecule 28: 50S ribosomal protein L35



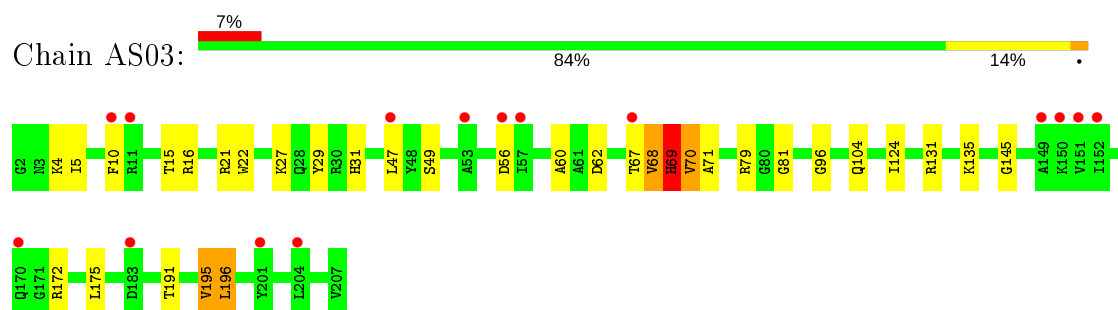
- Molecule 29: 30S ribosomal protein S2



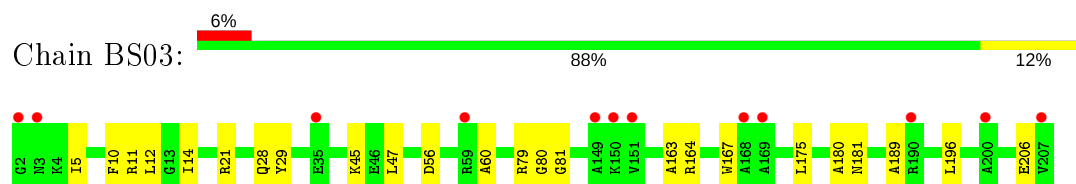
- Molecule 29: 30S ribosomal protein S2



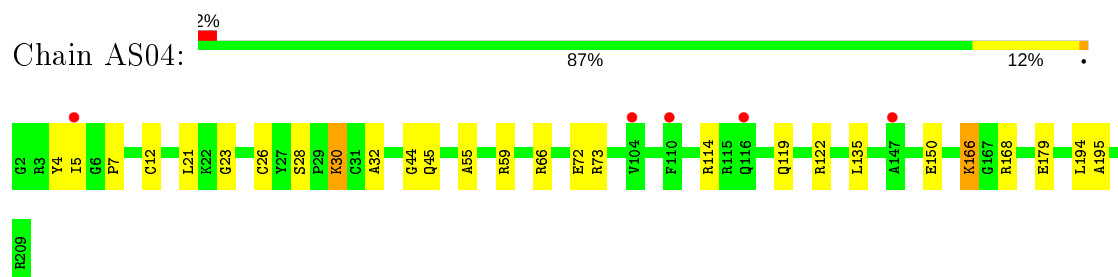
- Molecule 30: 30S ribosomal protein S3



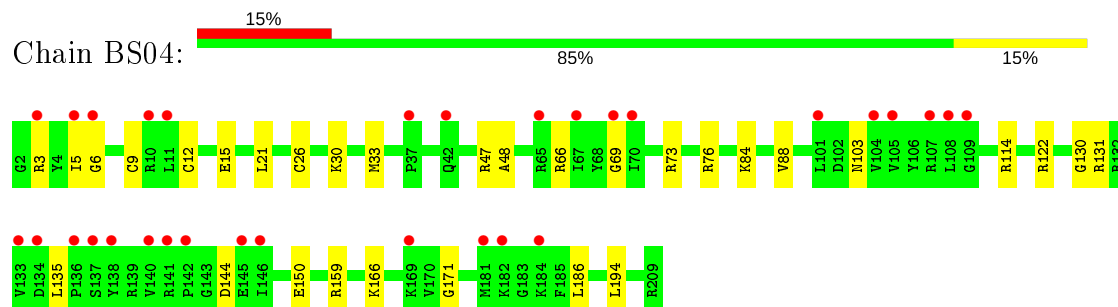
- Molecule 30: 30S ribosomal protein S3



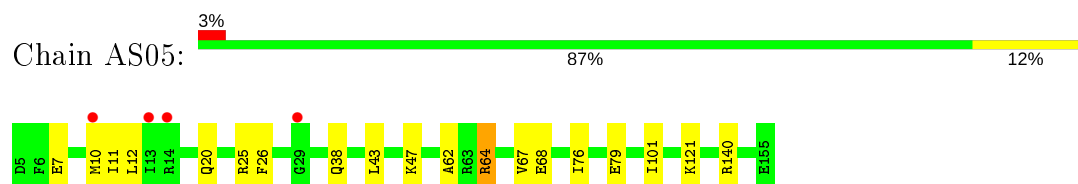
- Molecule 31: 30S ribosomal protein S4



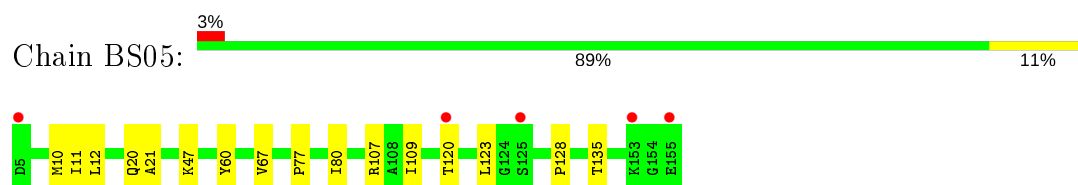
- Molecule 31: 30S ribosomal protein S4



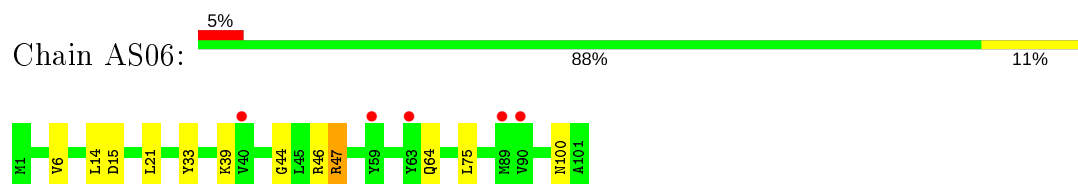
- Molecule 32: 30S ribosomal protein S5



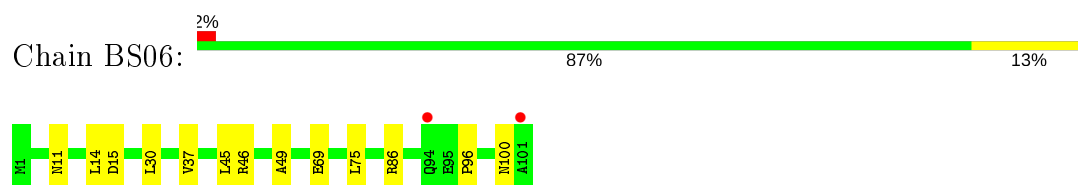
- Molecule 32: 30S ribosomal protein S5



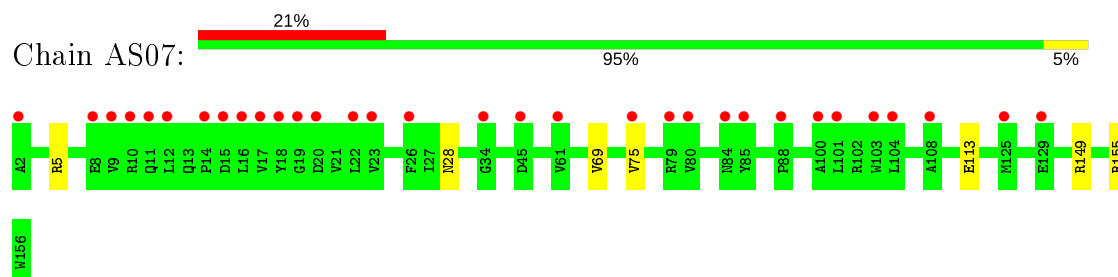
- Molecule 33: 30S ribosomal protein S6



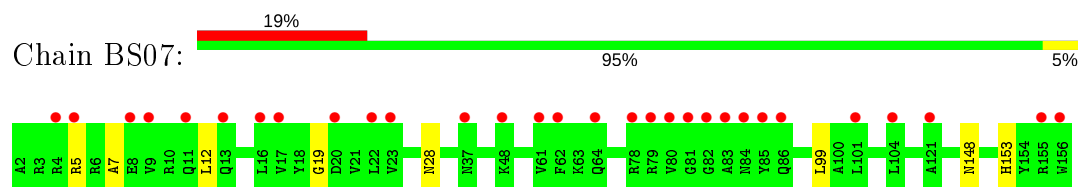
- Molecule 33: 30S ribosomal protein S6



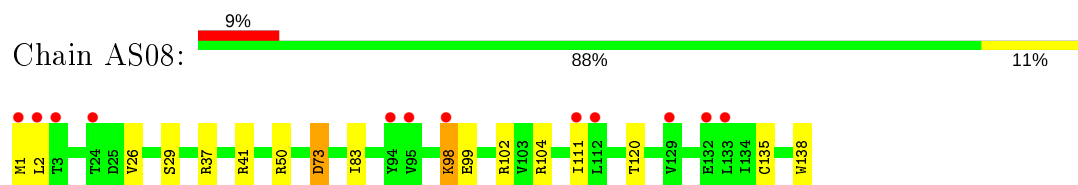
- Molecule 34: 30S ribosomal protein S7



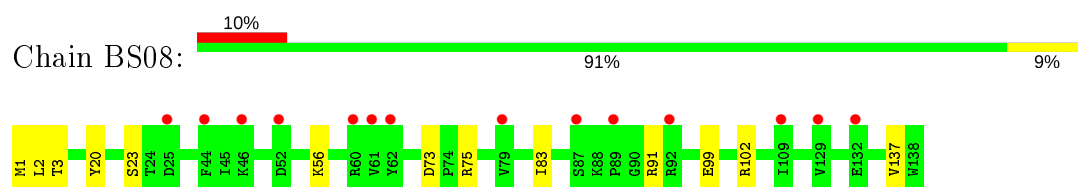
- Molecule 34: 30S ribosomal protein S7



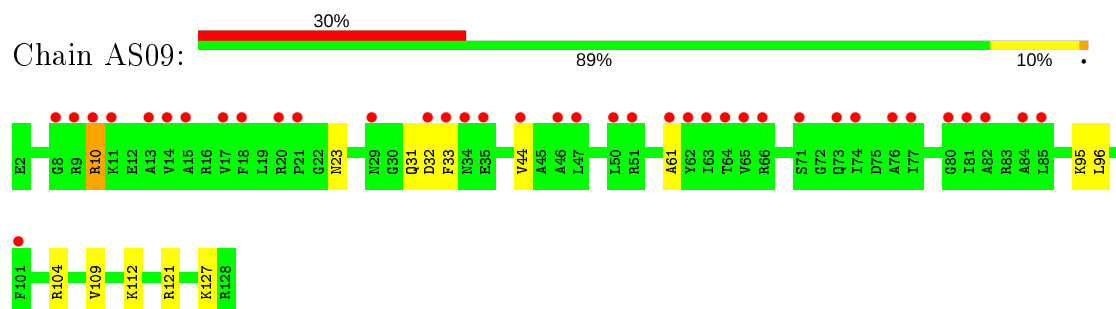
- Molecule 35: 30S ribosomal protein S8



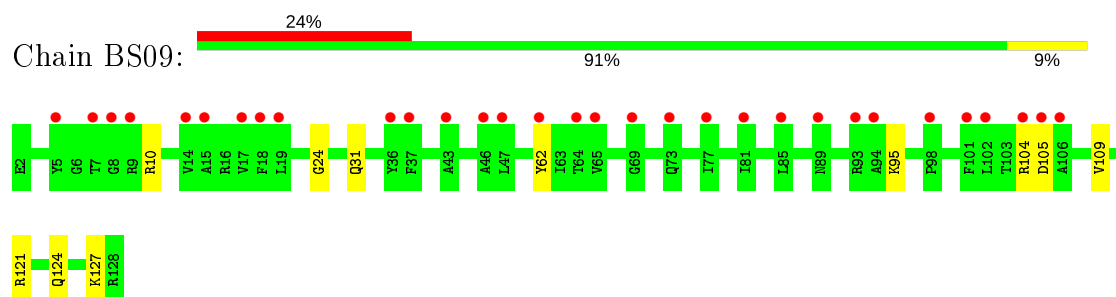
- Molecule 35: 30S ribosomal protein S8



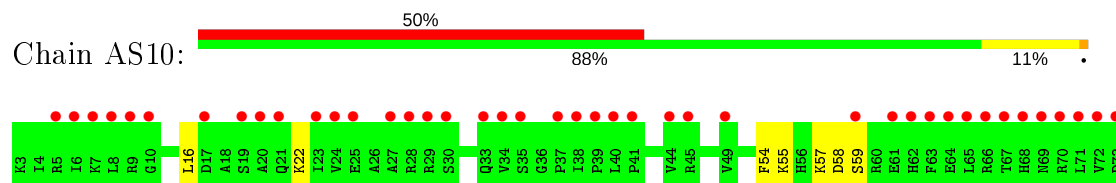
- Molecule 36: 30S ribosomal protein S9

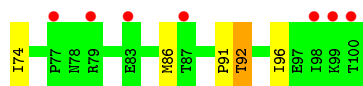


- Molecule 36: 30S ribosomal protein S9

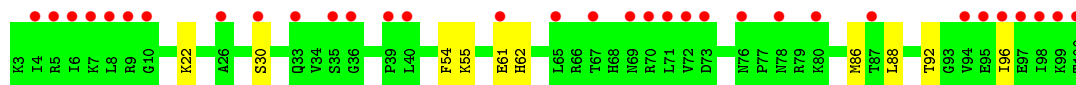


- Molecule 37: 30S ribosomal protein S10

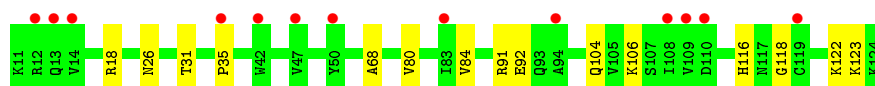
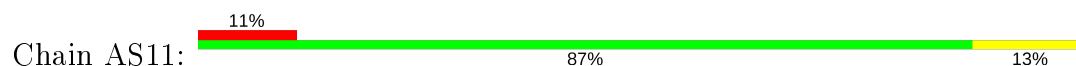




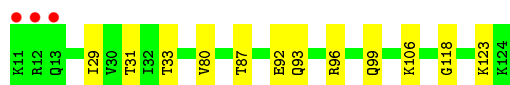
- Molecule 37: 30S ribosomal protein S10



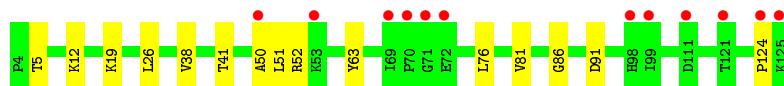
- Molecule 38: 30S ribosomal protein S11



- Molecule 38: 30S ribosomal protein S11



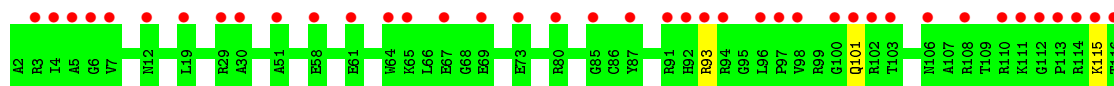
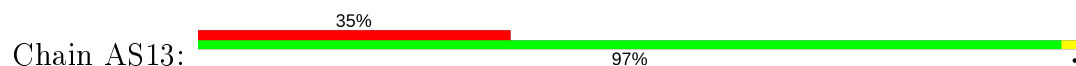
- Molecule 39: 30S ribosomal protein S12



- Molecule 39: 30S ribosomal protein S12

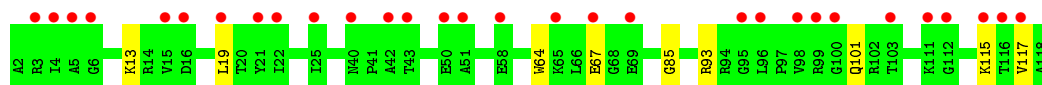


- Molecule 40: 30S ribosomal protein S13

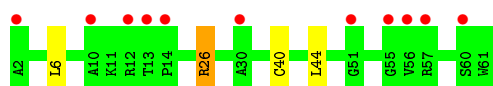




- Molecule 40: 30S ribosomal protein S13



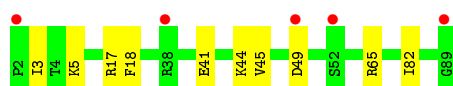
- Molecule 41: 30S ribosomal protein S14



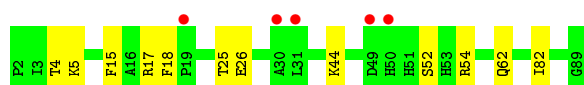
- Molecule 41: 30S ribosomal protein S14



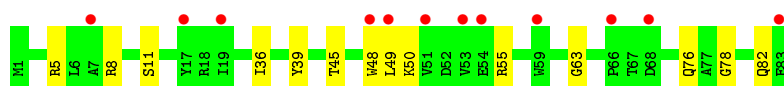
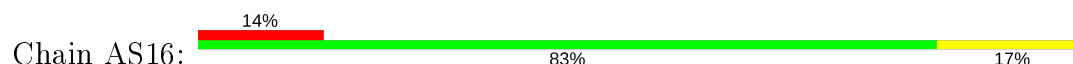
- Molecule 42: 30S ribosomal protein S15



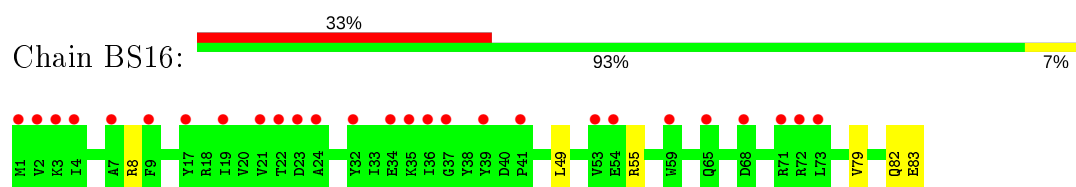
- Molecule 42: 30S ribosomal protein S15



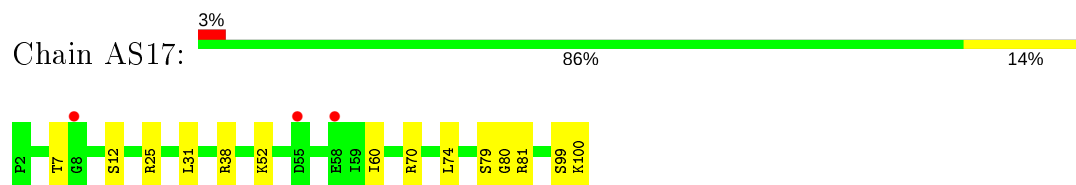
- Molecule 43: 30S ribosomal protein S16



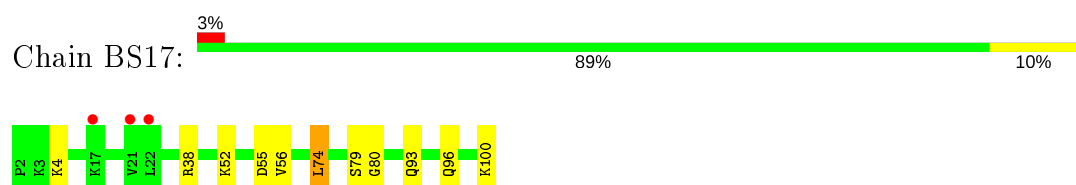
- Molecule 43: 30S ribosomal protein S16



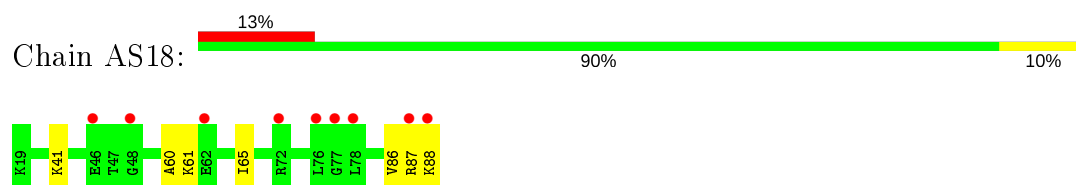
- Molecule 44: 30S ribosomal protein S17



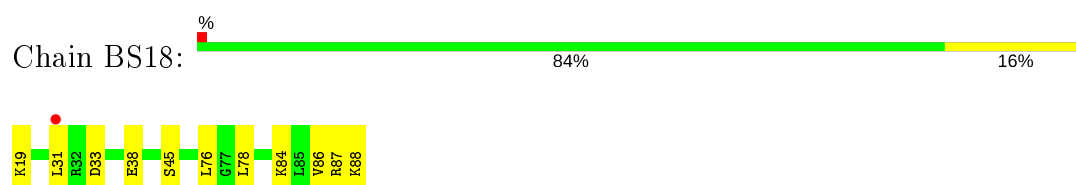
- Molecule 44: 30S ribosomal protein S17



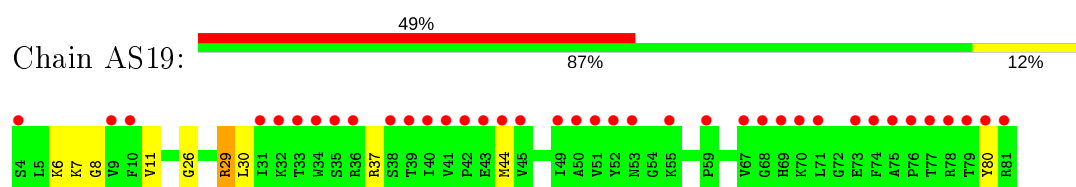
- Molecule 45: 30S ribosomal protein S18



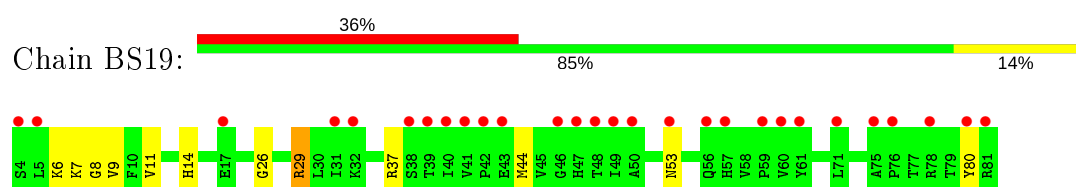
- Molecule 45: 30S ribosomal protein S18



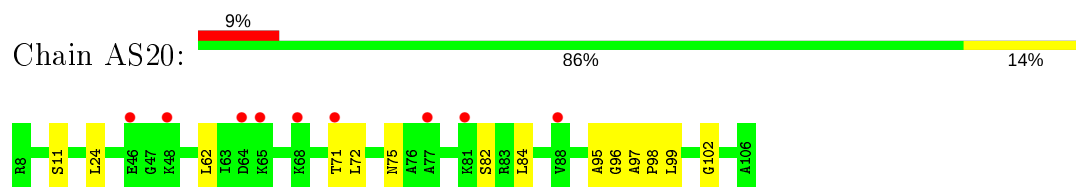
- Molecule 46: 30S ribosomal protein S19



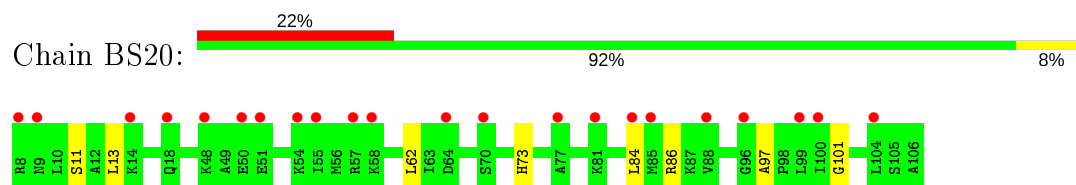
- Molecule 46: 30S ribosomal protein S19



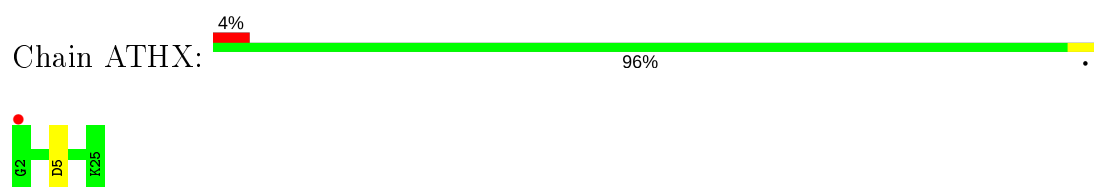
• Molecule 47: 30S ribosomal protein S20



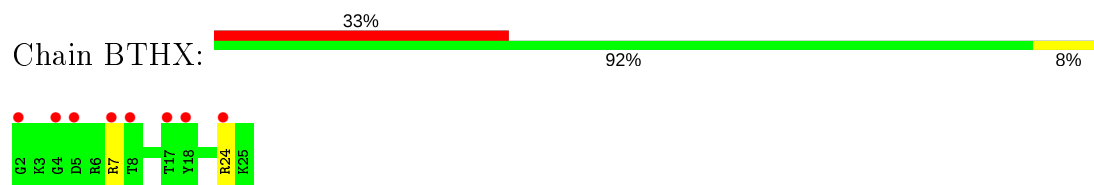
• Molecule 47: 30S ribosomal protein S20



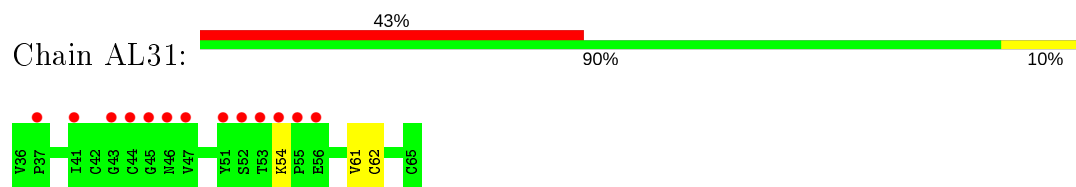
• Molecule 48: 30S ribosomal protein Thx



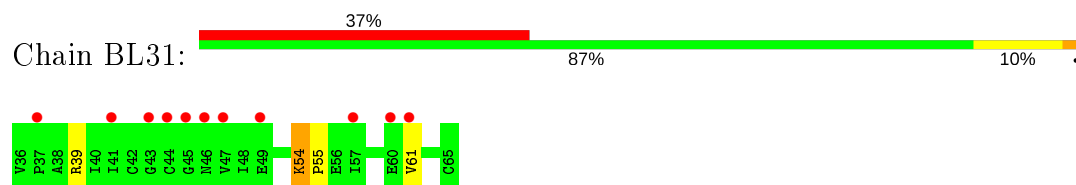
• Molecule 48: 30S ribosomal protein Thx



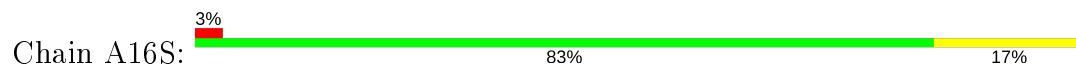
• Molecule 49: 50S ribosomal protein L31

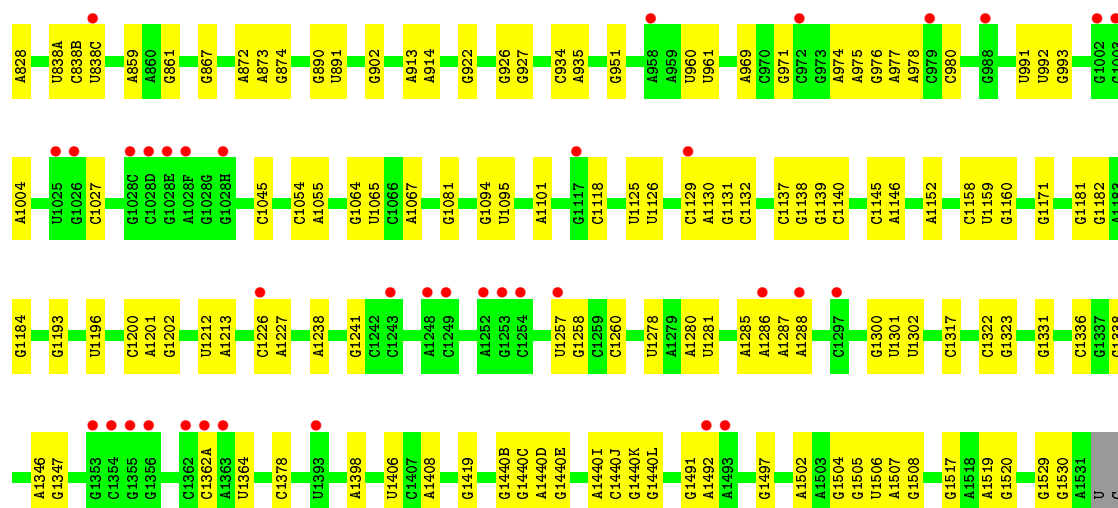


• Molecule 49: 50S ribosomal protein L31

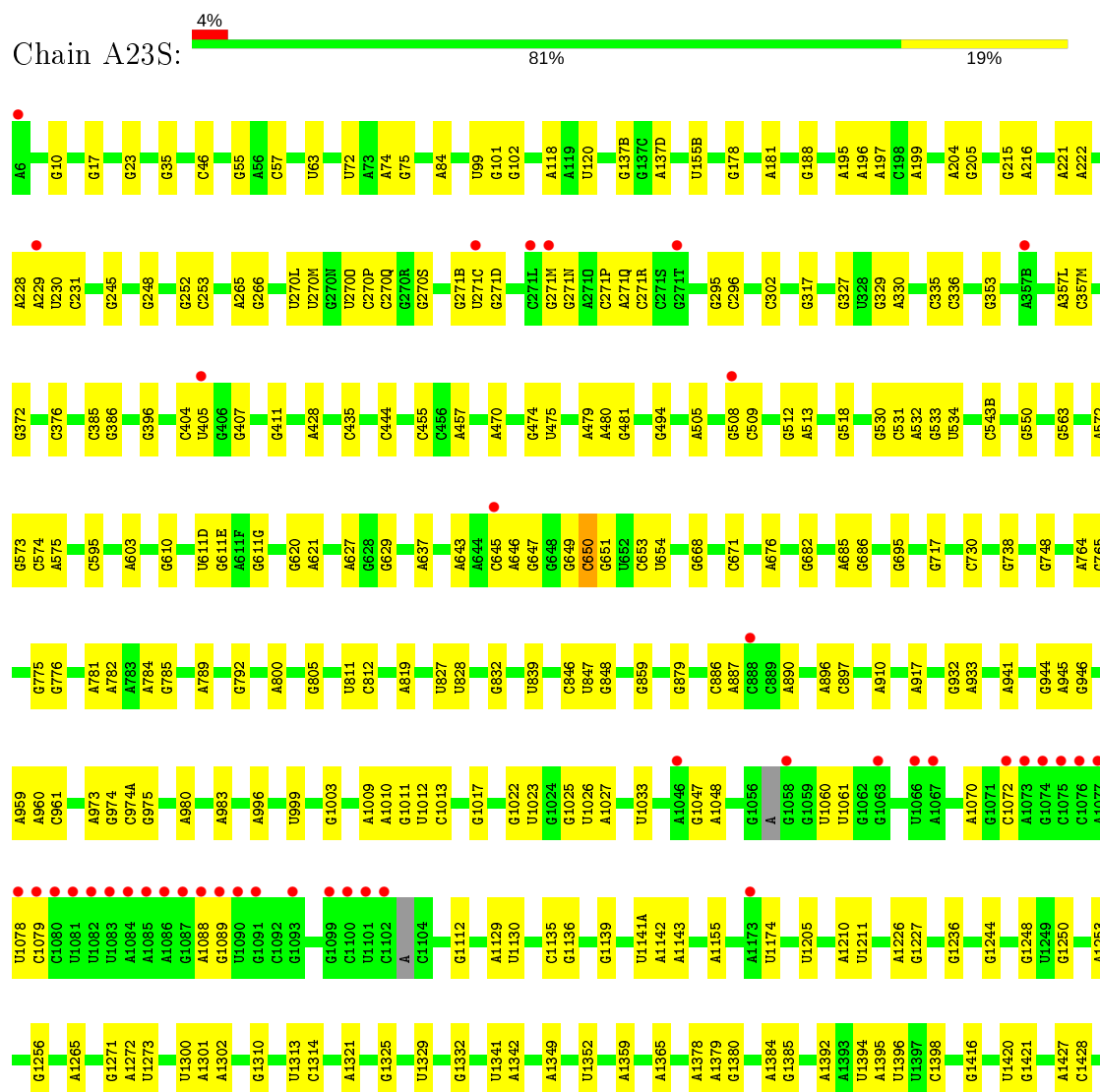


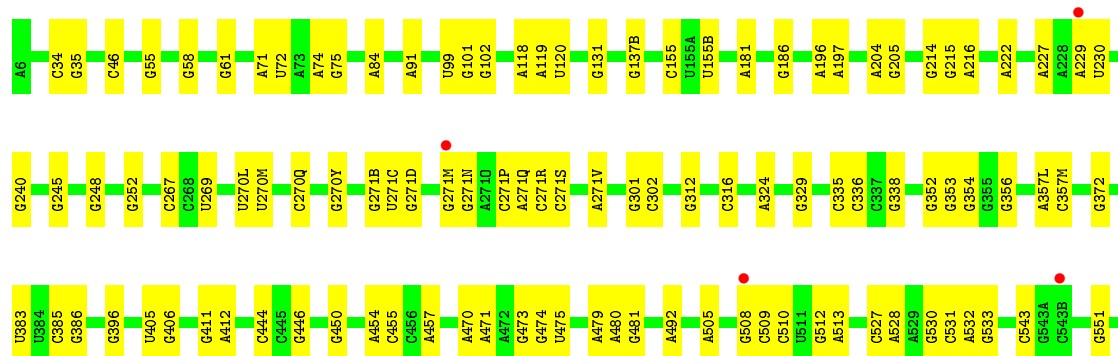
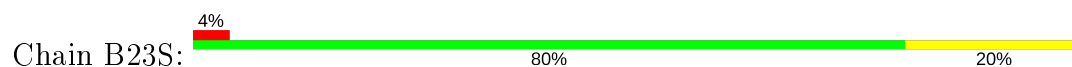
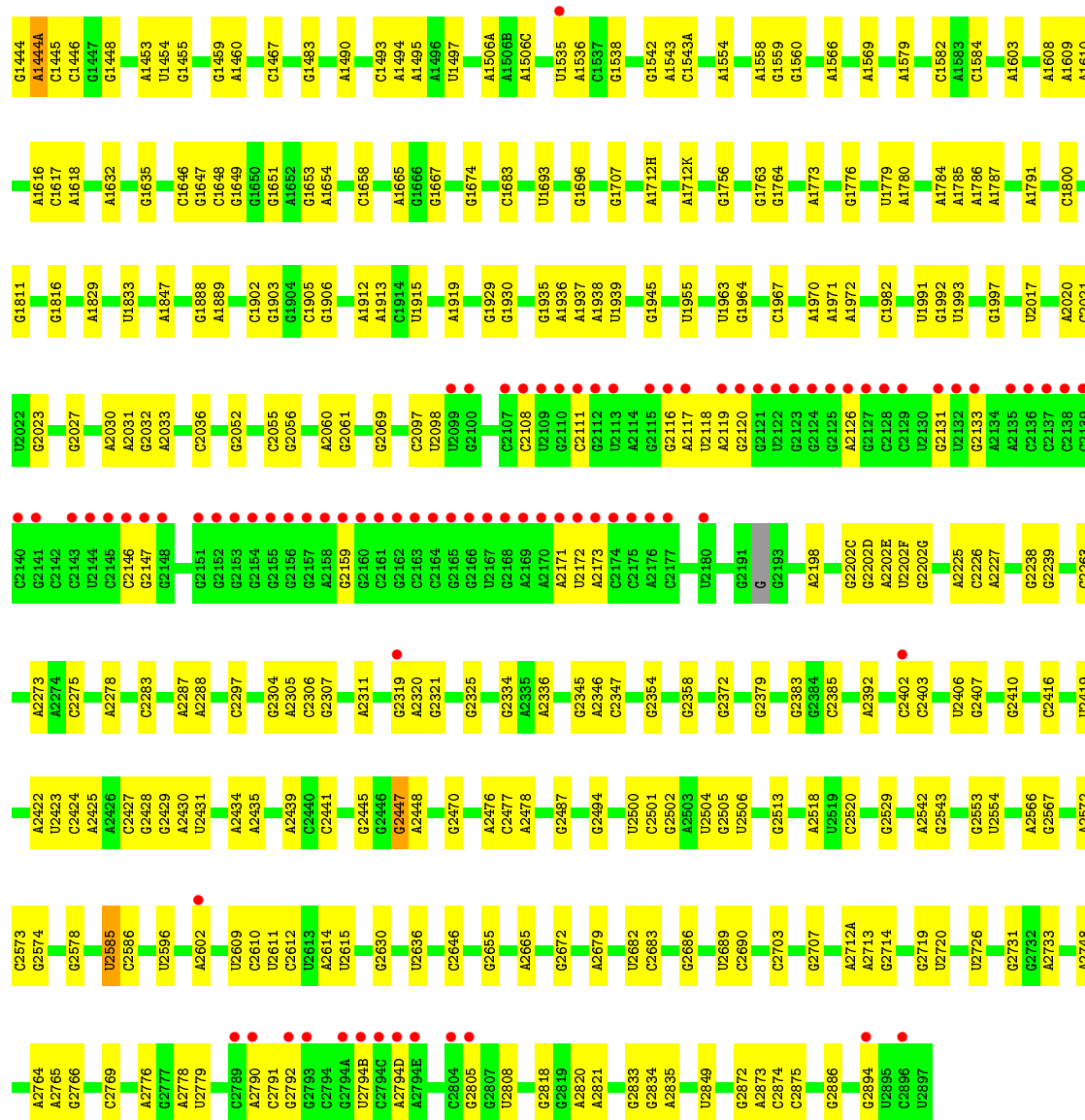
• Molecule 50: 16S ribosomal RNA



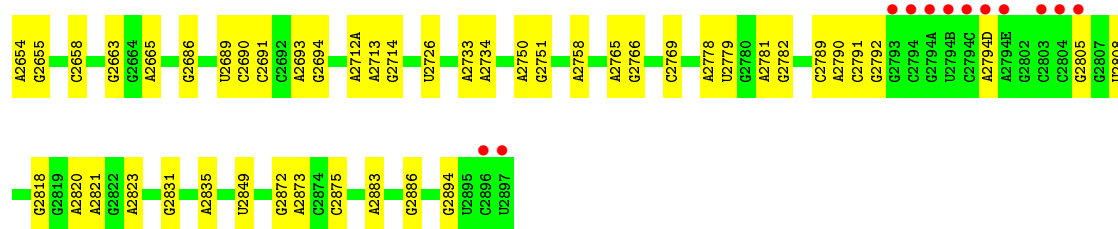


• Molecule 51: 23S ribosomal RNA

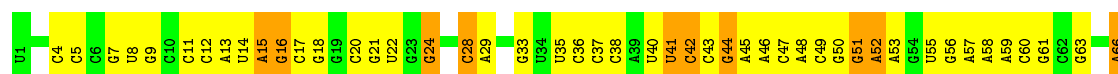




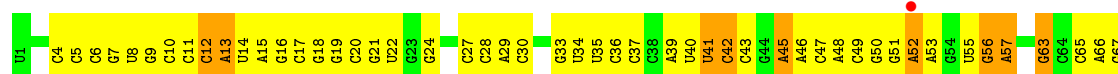
G2494	C2350	G2193	G2120	G1983	C1800	A1586	A1366	U1174	G1047	C914	A764	G563
U2500	A2361	A2198	G2121	G1984	G1801	A1603	A1379	G1175	A1048	C915	G765	G573
C2501	G2379	U1991	G2122	U1991	G1811	G1603	G1380	A1177	G1056	G916	G769	A574
G2502	G2202C	G1992	G2123	G1992	G1816	A1608	A1384	G1186	A	A918	A774	A575
A2503	G2202D	U1993	G2124	U1993	G1816	A1609	G1385	U1205	G1058	G919	G775	C587
U2504	A2202E		G2125		U1833	A1610	G1386	A1210	G1059	A926	G776	
G2505	U2202F	G1997	A2126	G1997	U1833	C1617		A1210	U1060			
U2506	G2202G	G2004	G2131	G2004	C1838			U1211	U1061	G932		G599
G2513	A2225	A2005	G2132	A2005	G1839	U1639	U1395	U1211	G1062	A782	A783	A603
A2518	C2226	C2006	G2133	C2006	G1840		U1416	A1219A	G1063	G938	A784	U611D
U2519	G2238	C2021	A2135	C2021	A1847	C1644		C1220	U1065		A785	G611E
C2520	G2239	U2022	G2136	U2022	G1858	G1645	A1427	G1227	U1066	A941	C786	G611F
G2529	U2243	G2023	C2137	G2023	G1858	G1646	C1428		G1067	A945	U787	U611G
U2542	U2406	A2031	C2138	A1859	G1859	C1648		G1236	G1068	G946	A788	
G2543	G2410	G2032	C2139	G1860	G1860		A1444A		A1069		A789	G620
U2554	G2411	A2033	G2141	G1878	G1878	G1651		U1240	A1070	A953	C790	A621
A2566	A2266	C2036	C2142			A1652		G1071	G1071		C791	A622
G2567	A2269	C2043	C2145	A1889	A1889	G1683	A1483	G1244	C1072	A959	G905	A627
G2568	A2273	G2043	G2146	G1903	G1903	G1674	U1484		G1073	A960		
G2569	A2274	G2052	G2147	G1904	G1904		G1485	G1248	G1074	C961	C812	A637
A2572	C2275	C2055	G2148	G1905	G1905	G1681	G1488	U1249	C1075	G974	A819	G645
U2573	A2278	A2057	G2149	G1906	G1906		G1489	G1250	C1076	C974A	A820	A646
A2578	C2283	A2058	U2150	G1913	G1913	G1685	A1480	G1252	U1077	G975	U827	G647
C2585	A2287	A2059	G2151	C1914	C1914	A1689	C1467	G1256	C1078	G979	U828	C650
U2586	U2296	G2061	G2152	A1919	A1919	C1694	G1483	G1271	U1081	A980	A829	G651
U2593	G2304	G2069	G2153	G1929	G1929	A1698	A1490	A1272	U1082	A983	G830	U652
G2602	A2305	G2080	G2154	C1934	C1934	A1712H	C1493		A1088	A990	G831	C653
U2603	G2306	C2081	G2155	G1935	G1935	U1712I	A1496	U1300	G1089	A996	U847	U654
U2604	G2307	U2086	G2156	A1937	A1937	G1712J	U1497	A1301	U1090		G848	G668
U2605	A2310		G2157	U1939	U1939	A1712K	A1506C	A1302	G1091	U999	A949	G669
U2609	G2317	U	G2158	G1945	G1945	G1712Q	U1535	G1309	C1092	G1003	G859	A670
C2610	G2318	U2099	G2159	U1955	U1955	G1750	G1538	U1343	A1098	C1004		G685
U2611	G2319	G2106	G2160	U1963	U1963	C1784	G1542	C1314	G1112	C1005	G879	G687
C2612	A2320	C2107	G2161	G1964	G1964	A1755	A1543	A1321	A1009	A1010	C384	U688
A2613	G2321	C2108	G2162	G1965	G1965	G1756	C1543A	U1329	G1011	G1011	C385	
U2614	G2325	U2109	G2163	A1966	A1966	G1763	A1558	C1330	U1012	U1012	A887	G717
U2615	G2334	C2110	G2164	C1967	C1967	G1764	G1559	A1331	C1013	C1013	C888	A722
C2616	A2335	G2111	A1773			A1773		G1332	C1135	A1020	C889	
G2630	A2336	U2112	A1785	A1970	A1970	A1785	A1566	U1344	G1136	A1021	A890	C730
C2634	U2344	A2114	A1786	A1971	A1971	A1786	A1569	C1345	U1141A	G1022	A896	G733
U2635	G2345	G2115	A1787	G1972	G1972	C1788	U1578	A1349	U1142	U1023	C897	G733
U2636	A2346	G2116	C1788	G1974	G1974		A1579	A1359	G1151	G1024	C898	U746
G2641	U2491	U2118	A1981	C1982	C1982	A1791	C1584	A1365	A1155	U1025	U907	G748
		A2119							A1173	U1026	A910	U762
			G							U1033		G763



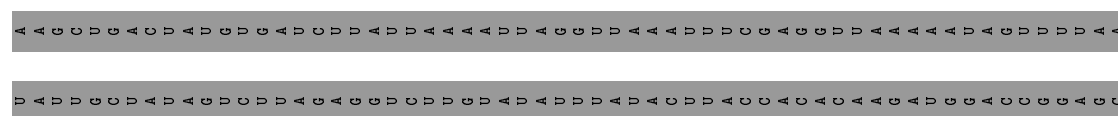
• Molecule 52: 5S ribosomal RNA



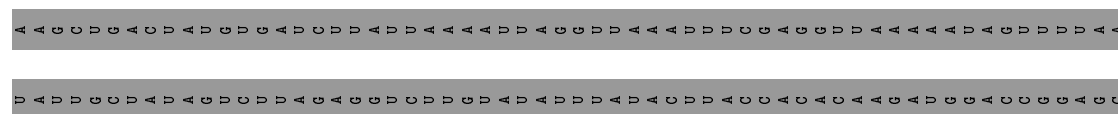
• Molecule 52: 5S ribosomal RNA



• Molecule 53: IRES RNA



• Molecule 53: IRES RNA





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	209.05Å 447.22Å 608.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.00 – 3.80 59.35 – 3.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (60.00-3.80) 99.9 (59.35-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 3.77Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.246 , 0.284 0.242 , 0.279	Depositor DCC
R_{free} test set	2000 reflections (0.36%)	wwPDB-VP
Wilson B-factor (Å ²)	119.0	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 93.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	287293	wwPDB-VP
Average B, all atoms (Å ²)	150.0	wwPDB-VP

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

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

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	AL02	0.33	0/2155	0.56	0/2905
1	BL02	0.40	0/2155	0.60	0/2905
2	AL03	0.31	0/1597	0.52	0/2153
2	BL03	0.33	0/1597	0.56	0/2153
3	AL04	0.31	0/1622	0.53	0/2194
3	BL04	0.32	0/1622	0.54	0/2194
4	AL05	0.22	0/1500	0.43	0/2017
4	BL05	0.23	0/1500	0.44	0/2017
5	AL06	0.25	0/1246	0.46	0/1682
5	BL06	0.33	0/1246	0.51	0/1682
6	AL09	0.29	0/1148	0.51	0/1552
6	BL09	0.30	0/1148	0.53	0/1552
7	AL11	0.22	0/1108	0.41	0/1500
7	BL11	0.25	0/1108	0.43	0/1500
8	AL13	0.37	0/1124	0.56	0/1515
8	BL13	0.32	0/1124	0.51	0/1515
9	AL14	0.29	0/942	0.51	0/1268
9	BL14	0.31	0/942	0.51	0/1268
10	AL15	0.31	0/1131	0.59	0/1504
10	BL15	0.37	0/1131	0.65	0/1504
11	AL16	0.31	0/1085	0.52	0/1449
11	BL16	0.32	0/1085	0.52	0/1449
12	AL17	0.37	0/974	0.57	0/1302
12	BL17	0.32	0/974	0.53	0/1302
13	AL18	0.25	0/779	0.43	0/1036
13	BL18	0.26	0/779	0.44	0/1036
14	AL19	0.27	0/1158	0.49	0/1544
14	BL19	0.27	0/1158	0.48	0/1544
15	AL20	0.32	0/982	0.47	0/1306
15	BL20	0.31	0/982	0.51	0/1306
16	AL21	0.31	0/790	0.52	0/1057
16	BL21	0.35	0/790	0.56	0/1057

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AL22	0.31	0/902	0.53	0/1209
17	BL22	0.32	0/902	0.51	0/1209
18	AL23	0.28	0/740	0.49	0/993
18	BL23	0.33	0/740	0.49	0/993
19	AL24	0.50	2/789 (0.3%)	1.34	4/1051 (0.4%)
19	BL24	0.34	0/789	0.58	0/1051
20	AL25	0.28	0/1515	0.46	0/2056
20	BL25	0.28	0/1515	0.48	0/2056
21	AL27	0.27	0/613	0.50	0/816
21	BL27	0.27	0/613	0.49	0/816
22	AL28	0.32	0/702	0.59	1/932 (0.1%)
22	BL28	0.34	0/702	0.60	0/932
23	AL29	0.27	0/523	0.49	0/690
23	BL29	0.32	0/523	0.55	0/690
24	AL30	0.24	0/473	0.44	0/634
24	BL30	0.27	0/473	0.47	0/634
25	AL32	0.38	0/419	0.59	0/567
25	BL32	0.41	0/419	0.63	0/567
26	AL33	0.28	0/388	0.52	0/518
26	BL33	0.24	0/388	0.47	0/518
27	AL34	0.35	0/427	0.49	0/561
27	BL34	0.39	0/427	0.54	0/561
28	AL35	0.36	0/516	0.56	0/679
28	BL35	0.43	0/516	0.70	1/679 (0.1%)
29	AS02	0.27	0/1936	0.46	1/2609 (0.0%)
29	BS02	0.26	0/1936	0.45	0/2609
30	AS03	0.89	8/1637 (0.5%)	0.95	13/2205 (0.6%)
30	BS03	0.26	0/1637	0.42	0/2205
31	AS04	0.30	0/1733	0.49	0/2318
31	BS04	0.25	0/1733	0.42	0/2318
32	AS05	0.24	0/1172	0.45	0/1576
32	BS05	0.25	0/1172	0.44	0/1576
33	AS06	0.23	0/856	0.44	0/1154
33	BS06	0.25	0/856	0.45	0/1154
34	AS07	0.25	0/1276	0.42	0/1709
34	BS07	0.26	0/1276	0.42	0/1709
35	AS08	0.23	0/1136	0.45	0/1527
35	BS08	0.24	0/1136	0.44	0/1527
36	AS09	0.25	0/1029	0.41	0/1378
36	BS09	0.21	0/1029	0.40	0/1378
37	AS10	0.28	0/808	0.47	0/1085
37	BS10	0.25	0/808	0.46	0/1085
38	AS11	0.27	0/857	0.48	0/1157

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	BS11	0.28	0/857	0.47	0/1157
39	AS12	0.29	0/973	0.50	0/1301
39	BS12	0.28	0/973	0.48	0/1301
40	AS13	0.27	0/944	0.43	0/1265
40	BS13	0.21	0/944	0.43	0/1265
41	AS14	0.30	0/501	0.45	0/664
41	BS14	0.22	0/501	0.39	0/664
42	AS15	0.39	0/745	0.54	0/992
42	BS15	0.38	0/745	0.52	0/992
43	AS16	0.31	0/717	0.48	0/963
43	BS16	0.25	0/717	0.46	0/963
44	AS17	0.31	0/837	0.48	0/1117
44	BS17	0.30	0/837	0.49	0/1117
45	AS18	0.24	0/579	0.43	0/768
45	BS18	0.25	0/579	0.45	0/768
46	AS19	0.30	0/643	0.46	0/865
46	BS19	0.27	0/643	0.46	0/865
47	AS20	0.26	0/764	0.47	0/1006
47	BS20	0.27	0/764	0.47	0/1006
48	ATHX	0.19	0/213	0.37	0/277
48	BTHX	0.20	0/213	0.40	0/277
49	AL31	0.20	0/229	0.45	0/309
49	BL31	0.20	0/229	0.41	0/309
50	A16S	0.27	0/36194	0.59	0/56493
50	B16S	0.25	0/36193	0.57	0/56490
51	A23S	0.39	2/69356 (0.0%)	0.71	11/108266 (0.0%)
51	B23S	0.39	0/69359	0.74	23/108270 (0.0%)
52	A5S	0.24	0/2853	0.55	0/4451
52	B5S	0.24	0/2853	0.56	0/4451
53	AIRE	0.25	0/748	0.50	0/1160
53	BIRE	0.27	0/748	0.57	0/1160
All	All	0.34	12/312170 (0.0%)	0.64	54/466611 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
10	AL15	0	1
10	BL15	0	1
19	AL24	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
30	AS03	0	2
All	All	0	5

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
51	A23S	253	C	C4-N4	40.77	1.70	1.33
51	A23S	253	C	N3-C4	19.94	1.48	1.33
30	AS03	69	HIS	C-O	13.83	1.49	1.23
30	AS03	70	VAL	CB-CG1	13.29	1.80	1.52
30	AS03	69	HIS	N-CA	12.82	1.72	1.46

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	A23S	1444(A)	A	P-O3'-C3'	37.85	165.12	119.70
19	AL24	6	HIS	CG-ND1-CE1	-32.52	62.67	108.20
30	AS03	69	HIS	CB-CG-ND1	-19.53	74.38	123.20
19	AL24	6	HIS	CG-CD2-NE2	-17.32	76.28	109.20
30	AS03	69	HIS	N-CA-CB	17.00	141.21	110.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
10	AL15	51	PHE	Peptide
19	AL24	6	HIS	Sidechain
30	AS03	68	VAL	Mainchain
30	AS03	69	HIS	Sidechain
10	BL15	51	PHE	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	AL02	2105	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BL02	2105	0	0	0	0
2	AL03	1564	0	0	0	0
2	BL03	1564	0	0	0	0
3	AL04	1587	0	0	0	0
3	BL04	1587	0	0	0	0
4	AL05	1475	0	0	0	0
4	BL05	1475	0	0	0	0
5	AL06	1223	0	0	0	0
5	BL06	1223	0	0	0	0
6	AL09	1133	0	0	0	0
6	BL09	1133	0	0	0	0
7	AL11	1088	0	0	0	0
7	BL11	1088	0	0	0	0
8	AL13	1097	0	0	0	0
8	BL13	1097	0	0	0	0
9	AL14	932	0	0	0	0
9	BL14	932	0	0	0	0
10	AL15	1114	0	0	0	0
10	BL15	1114	0	0	0	0
11	AL16	1065	0	0	0	0
11	BL16	1065	0	0	0	0
12	AL17	960	0	0	0	0
12	BL17	960	0	0	0	0
13	AL18	771	0	0	0	0
13	BL18	771	0	0	0	0
14	AL19	1144	0	0	0	0
14	BL19	1144	0	0	0	0
15	AL20	964	0	0	0	0
15	BL20	964	0	0	0	0
16	AL21	779	0	0	0	0
16	BL21	779	0	0	0	0
17	AL22	891	0	0	0	0
17	BL22	891	0	0	0	0
18	AL23	726	0	0	0	0
18	BL23	726	0	0	0	0
19	AL24	776	0	0	0	0
19	BL24	776	0	0	0	0
20	AL25	1483	0	0	0	0
20	BL25	1483	0	0	0	0
21	AL27	605	0	0	0	0
21	BL27	605	0	0	0	0
22	AL28	695	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
22	BL28	695	0	0	0	0
23	AL29	521	0	0	0	0
23	BL29	521	0	0	0	0
24	AL30	468	0	0	0	0
24	BL30	468	0	0	0	0
25	AL32	405	0	0	0	0
25	BL32	405	0	0	0	0
26	AL33	381	0	0	0	0
26	BL33	381	0	0	0	0
27	AL34	419	0	0	0	0
27	BL34	419	0	0	0	0
28	AL35	508	0	0	0	0
28	BL35	508	0	0	0	0
29	AS02	1901	0	0	0	0
29	BS02	1901	0	0	0	0
30	AS03	1613	0	0	0	0
30	BS03	1613	0	0	0	0
31	AS04	1703	0	0	0	0
31	BS04	1703	0	0	0	0
32	AS05	1156	0	0	0	0
32	BS05	1156	0	0	0	0
33	AS06	843	0	0	0	0
33	BS06	843	0	0	0	0
34	AS07	1257	0	0	0	0
34	BS07	1257	0	0	0	0
35	AS08	1116	0	0	0	0
35	BS08	1116	0	0	0	0
36	AS09	1011	0	0	0	0
36	BS09	1011	0	0	0	0
37	AS10	795	0	0	0	0
37	BS10	795	0	0	0	0
38	AS11	843	0	0	0	0
38	BS11	843	0	0	0	0
39	AS12	957	0	0	0	0
39	BS12	957	0	0	0	0
40	AS13	934	0	0	0	0
40	BS13	934	0	0	0	0
41	AS14	492	0	0	0	0
41	BS14	492	0	0	0	0
42	AS15	734	0	0	0	0
42	BS15	734	0	0	0	0
43	AS16	701	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
43	BS16	701	0	0	0	0
44	AS17	824	0	0	0	0
44	BS17	824	0	0	0	0
45	AS18	574	0	0	0	0
45	BS18	574	0	0	0	0
46	AS19	630	0	0	0	0
46	BS19	630	0	0	0	0
47	AS20	762	0	0	0	0
47	BS20	762	0	0	0	0
48	ATHX	209	0	0	0	0
48	BTHX	209	0	0	0	0
49	AL31	226	0	0	0	0
49	BL31	226	0	0	0	0
50	A16S	32332	0	0	0	0
50	B16S	32331	0	0	0	0
51	A23S	61929	0	0	0	0
51	B23S	61931	0	0	0	0
52	A5S	2551	0	1295	97	0
52	B5S	2551	0	1295	104	0
53	AIRE	672	0	0	0	0
53	BIRE	672	0	0	0	0
54	AS04	1	0	0	0	0
54	AS14	1	0	0	0	0
54	BS04	1	0	0	0	0
54	BS14	1	0	0	0	0
All	All	287293	0	2590	201	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 34.

The worst 5 of 201 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)
52:A5S:51:G:H21	52:A5S:52:A:H62	1.05	0.95
52:B5S:40:U:H3'	52:B5S:41:U:H5''	1.50	0.89
52:A5S:40:U:H3'	52:A5S:41:U:H5''	1.55	0.89
52:B5S:10:C:C2	52:B5S:11:C:H5	1.90	0.88
52:A5S:56:G:H21	52:A5S:59:A:H61	1.21	0.86

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AL02	269/271 (99%)	197 (73%)	49 (18%)	23 (9%)	1	12
1	BL02	269/271 (99%)	184 (68%)	56 (21%)	29 (11%)	0	8
2	AL03	202/204 (99%)	145 (72%)	41 (20%)	16 (8%)	1	14
2	BL03	202/204 (99%)	140 (69%)	46 (23%)	16 (8%)	1	14
3	AL04	200/202 (99%)	154 (77%)	31 (16%)	15 (8%)	1	15
3	BL04	200/202 (99%)	142 (71%)	41 (20%)	17 (8%)	1	12
4	AL05	179/181 (99%)	134 (75%)	35 (20%)	10 (6%)	2	21
4	BL05	179/181 (99%)	128 (72%)	35 (20%)	16 (9%)	1	12
5	AL06	157/159 (99%)	117 (74%)	35 (22%)	5 (3%)	4	32
5	BL06	157/159 (99%)	119 (76%)	27 (17%)	11 (7%)	1	17
6	AL09	143/145 (99%)	112 (78%)	26 (18%)	5 (4%)	3	31
6	BL09	143/145 (99%)	97 (68%)	37 (26%)	9 (6%)	1	20
7	AL11	145/147 (99%)	100 (69%)	36 (25%)	9 (6%)	1	20
7	BL11	145/147 (99%)	108 (74%)	28 (19%)	9 (6%)	1	20
8	AL13	135/137 (98%)	95 (70%)	23 (17%)	17 (13%)	0	5
8	BL13	135/137 (98%)	90 (67%)	30 (22%)	15 (11%)	0	7
9	AL14	120/122 (98%)	95 (79%)	16 (13%)	9 (8%)	1	15
9	BL14	120/122 (98%)	94 (78%)	15 (12%)	11 (9%)	1	12
10	AL15	144/146 (99%)	76 (53%)	36 (25%)	32 (22%)	0	1
10	BL15	144/146 (99%)	83 (58%)	41 (28%)	20 (14%)	0	4
11	AL16	132/134 (98%)	93 (70%)	28 (21%)	11 (8%)	1	13
11	BL16	132/134 (98%)	99 (75%)	20 (15%)	13 (10%)	0	10
12	AL17	115/117 (98%)	89 (77%)	17 (15%)	9 (8%)	1	15
12	BL17	115/117 (98%)	82 (71%)	27 (24%)	6 (5%)	2	23
13	AL18	96/98 (98%)	66 (69%)	17 (18%)	13 (14%)	0	4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	BL18	96/98 (98%)	61 (64%)	24 (25%)	11 (12%)	0	7
14	AL19	135/137 (98%)	85 (63%)	32 (24%)	18 (13%)	0	4
14	BL19	135/137 (98%)	88 (65%)	31 (23%)	16 (12%)	0	6
15	AL20	115/117 (98%)	88 (76%)	21 (18%)	6 (5%)	2	23
15	BL20	115/117 (98%)	88 (76%)	22 (19%)	5 (4%)	2	26
16	AL21	99/101 (98%)	70 (71%)	23 (23%)	6 (6%)	1	20
16	BL21	99/101 (98%)	75 (76%)	14 (14%)	10 (10%)	0	9
17	AL22	110/112 (98%)	86 (78%)	19 (17%)	5 (4%)	2	25
17	BL22	110/112 (98%)	81 (74%)	24 (22%)	5 (4%)	2	25
18	AL23	90/92 (98%)	78 (87%)	10 (11%)	2 (2%)	6	39
18	BL23	90/92 (98%)	78 (87%)	10 (11%)	2 (2%)	6	39
19	AL24	98/100 (98%)	62 (63%)	23 (24%)	13 (13%)	0	4
19	BL24	98/100 (98%)	53 (54%)	28 (29%)	17 (17%)	0	3
20	AL25	185/187 (99%)	144 (78%)	34 (18%)	7 (4%)	3	29
20	BL25	185/187 (99%)	142 (77%)	37 (20%)	6 (3%)	4	32
21	AL27	74/76 (97%)	56 (76%)	12 (16%)	6 (8%)	1	14
21	BL27	74/76 (97%)	58 (78%)	12 (16%)	4 (5%)	2	22
22	AL28	86/88 (98%)	60 (70%)	12 (14%)	14 (16%)	0	3
22	BL28	86/88 (98%)	58 (67%)	14 (16%)	14 (16%)	0	3
23	AL29	60/62 (97%)	45 (75%)	12 (20%)	3 (5%)	2	23
23	BL29	60/62 (97%)	44 (73%)	9 (15%)	7 (12%)	0	6
24	AL30	57/59 (97%)	49 (86%)	6 (10%)	2 (4%)	3	31
24	BL30	57/59 (97%)	41 (72%)	13 (23%)	3 (5%)	2	23
25	AL32	50/52 (96%)	36 (72%)	10 (20%)	4 (8%)	1	14
25	BL32	50/52 (96%)	40 (80%)	7 (14%)	3 (6%)	1	20
26	AL33	42/44 (96%)	27 (64%)	12 (29%)	3 (7%)	1	17
26	BL33	42/44 (96%)	26 (62%)	12 (29%)	4 (10%)	0	11
27	AL34	46/48 (96%)	40 (87%)	4 (9%)	2 (4%)	2	26
27	BL34	46/48 (96%)	36 (78%)	7 (15%)	3 (6%)	1	19
28	AL35	61/63 (97%)	36 (59%)	20 (33%)	5 (8%)	1	13
28	BL35	61/63 (97%)	39 (64%)	14 (23%)	8 (13%)	0	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
29	AS02	232/234 (99%)	184 (79%)	34 (15%)	14 (6%)	1	20
29	BS02	232/234 (99%)	183 (79%)	44 (19%)	5 (2%)	6	39
30	AS03	204/206 (99%)	159 (78%)	32 (16%)	13 (6%)	1	20
30	BS03	204/206 (99%)	146 (72%)	47 (23%)	11 (5%)	2	22
31	AS04	206/208 (99%)	154 (75%)	40 (19%)	12 (6%)	1	21
31	BS04	206/208 (99%)	153 (74%)	40 (19%)	13 (6%)	1	20
32	AS05	149/151 (99%)	114 (76%)	30 (20%)	5 (3%)	3	31
32	BS05	149/151 (99%)	120 (80%)	23 (15%)	6 (4%)	3	28
33	AS06	99/101 (98%)	80 (81%)	14 (14%)	5 (5%)	2	23
33	BS06	99/101 (98%)	79 (80%)	14 (14%)	6 (6%)	1	20
34	AS07	153/155 (99%)	132 (86%)	19 (12%)	2 (1%)	12	48
34	BS07	153/155 (99%)	128 (84%)	22 (14%)	3 (2%)	7	41
35	AS08	136/138 (99%)	113 (83%)	15 (11%)	8 (6%)	1	21
35	BS08	136/138 (99%)	106 (78%)	24 (18%)	6 (4%)	2	25
36	AS09	125/127 (98%)	94 (75%)	23 (18%)	8 (6%)	1	20
36	BS09	125/127 (98%)	98 (78%)	23 (18%)	4 (3%)	4	32
37	AS10	96/98 (98%)	76 (79%)	16 (17%)	4 (4%)	3	26
37	BS10	96/98 (98%)	73 (76%)	21 (22%)	2 (2%)	7	40
38	AS11	112/114 (98%)	85 (76%)	19 (17%)	8 (7%)	1	17
38	BS11	112/114 (98%)	91 (81%)	18 (16%)	3 (3%)	5	35
39	AS12	120/122 (98%)	90 (75%)	26 (22%)	4 (3%)	4	32
39	BS12	120/122 (98%)	91 (76%)	25 (21%)	4 (3%)	4	32
40	AS13	115/117 (98%)	96 (84%)	17 (15%)	2 (2%)	9	43
40	BS13	115/117 (98%)	93 (81%)	17 (15%)	5 (4%)	2	26
41	AS14	58/60 (97%)	47 (81%)	10 (17%)	1 (2%)	9	43
41	BS14	58/60 (97%)	47 (81%)	9 (16%)	2 (3%)	3	31
42	AS15	86/88 (98%)	69 (80%)	15 (17%)	2 (2%)	6	38
42	BS15	86/88 (98%)	69 (80%)	16 (19%)	1 (1%)	13	50
43	AS16	81/83 (98%)	61 (75%)	15 (18%)	5 (6%)	1	20
43	BS16	81/83 (98%)	63 (78%)	16 (20%)	2 (2%)	5	36
44	AS17	97/99 (98%)	74 (76%)	18 (19%)	5 (5%)	2	23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
44	BS17	97/99 (98%)	81 (84%)	11 (11%)	5 (5%)	2	23
45	AS18	68/70 (97%)	48 (71%)	16 (24%)	4 (6%)	1	21
45	BS18	68/70 (97%)	45 (66%)	20 (29%)	3 (4%)	2	25
46	AS19	76/78 (97%)	53 (70%)	18 (24%)	5 (7%)	1	19
46	BS19	76/78 (97%)	50 (66%)	20 (26%)	6 (8%)	1	14
47	AS20	97/99 (98%)	73 (75%)	14 (14%)	10 (10%)	0	9
47	BS20	97/99 (98%)	72 (74%)	21 (22%)	4 (4%)	3	27
48	ATHX	22/24 (92%)	18 (82%)	4 (18%)	0	100	100
48	BTHX	22/24 (92%)	13 (59%)	7 (32%)	2 (9%)	1	12
49	AL31	28/30 (93%)	15 (54%)	11 (39%)	2 (7%)	1	17
49	BL31	28/30 (93%)	20 (71%)	5 (18%)	3 (11%)	0	8
All	All	11410/11606 (98%)	8465 (74%)	2170 (19%)	775 (7%)	1	18

5 of 775 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AL02	26	LYS
1	AL02	33	LEU
1	AL02	34	VAL
1	AL02	154	LYS
1	AL02	239	ARG

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	
1	AL02	213/213 (100%)	185 (87%)	28 (13%)	4	23
1	BL02	213/213 (100%)	177 (83%)	36 (17%)	2	14
2	AL03	165/165 (100%)	137 (83%)	28 (17%)	2	14
2	BL03	165/165 (100%)	131 (79%)	34 (21%)	1	8
3	AL04	161/161 (100%)	150 (93%)	11 (7%)	16	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	BL04	161/161 (100%)	141 (88%)	20 (12%)	4	24
4	AL05	155/155 (100%)	133 (86%)	22 (14%)	3	21
4	BL05	155/155 (100%)	136 (88%)	19 (12%)	4	24
5	AL06	132/132 (100%)	123 (93%)	9 (7%)	16	47
5	BL06	132/132 (100%)	116 (88%)	16 (12%)	5	25
6	AL09	122/122 (100%)	97 (80%)	25 (20%)	1	8
6	BL09	122/122 (100%)	102 (84%)	20 (16%)	2	15
7	AL11	111/111 (100%)	105 (95%)	6 (5%)	22	53
7	BL11	111/111 (100%)	106 (96%)	5 (4%)	27	57
8	AL13	116/116 (100%)	100 (86%)	16 (14%)	3	22
8	BL13	116/116 (100%)	98 (84%)	18 (16%)	2	17
9	AL14	100/100 (100%)	87 (87%)	13 (13%)	4	23
9	BL14	100/100 (100%)	89 (89%)	11 (11%)	6	29
10	AL15	112/112 (100%)	82 (73%)	30 (27%)	0	3
10	BL15	112/112 (100%)	80 (71%)	32 (29%)	0	2
11	AL16	105/105 (100%)	96 (91%)	9 (9%)	10	40
11	BL16	105/105 (100%)	88 (84%)	17 (16%)	2	15
12	AL17	100/100 (100%)	87 (87%)	13 (13%)	4	23
12	BL17	100/100 (100%)	86 (86%)	14 (14%)	3	21
13	AL18	77/77 (100%)	70 (91%)	7 (9%)	9	36
13	BL18	77/77 (100%)	68 (88%)	9 (12%)	5	27
14	AL19	121/121 (100%)	105 (87%)	16 (13%)	4	22
14	BL19	121/121 (100%)	106 (88%)	15 (12%)	4	24
15	AL20	93/93 (100%)	79 (85%)	14 (15%)	3	18
15	BL20	93/93 (100%)	83 (89%)	10 (11%)	6	30
16	AL21	82/82 (100%)	68 (83%)	14 (17%)	2	14
16	BL21	82/82 (100%)	67 (82%)	15 (18%)	1	11
17	AL22	91/91 (100%)	83 (91%)	8 (9%)	10	38
17	BL22	91/91 (100%)	83 (91%)	8 (9%)	10	38
18	AL23	74/74 (100%)	67 (90%)	7 (10%)	8	34
18	BL23	74/74 (100%)	67 (90%)	7 (10%)	8	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
19	AL24	84/84 (100%)	69 (82%)	15 (18%)	2	12
19	BL24	84/84 (100%)	71 (84%)	13 (16%)	2	17
20	AL25	162/162 (100%)	150 (93%)	12 (7%)	13	44
20	BL25	162/162 (100%)	151 (93%)	11 (7%)	16	47
21	AL27	61/61 (100%)	56 (92%)	5 (8%)	11	40
21	BL27	61/61 (100%)	52 (85%)	9 (15%)	3	19
22	AL28	73/73 (100%)	60 (82%)	13 (18%)	2	12
22	BL28	73/73 (100%)	60 (82%)	13 (18%)	2	12
23	AL29	58/58 (100%)	52 (90%)	6 (10%)	7	31
23	BL29	58/58 (100%)	47 (81%)	11 (19%)	1	10
24	AL30	51/51 (100%)	47 (92%)	4 (8%)	12	42
24	BL30	51/51 (100%)	42 (82%)	9 (18%)	2	13
25	AL32	45/45 (100%)	38 (84%)	7 (16%)	2	17
25	BL32	45/45 (100%)	40 (89%)	5 (11%)	6	29
26	AL33	43/43 (100%)	36 (84%)	7 (16%)	2	15
26	BL33	43/43 (100%)	34 (79%)	9 (21%)	1	8
27	AL34	41/41 (100%)	37 (90%)	4 (10%)	8	33
27	BL34	41/41 (100%)	33 (80%)	8 (20%)	1	9
28	AL35	53/53 (100%)	45 (85%)	8 (15%)	3	18
28	BL35	53/53 (100%)	49 (92%)	4 (8%)	13	44
29	AS02	202/202 (100%)	186 (92%)	16 (8%)	12	42
29	BS02	202/202 (100%)	186 (92%)	16 (8%)	12	42
30	AS03	160/160 (100%)	140 (88%)	20 (12%)	4	24
30	BS03	160/160 (100%)	147 (92%)	13 (8%)	11	41
31	AS04	180/180 (100%)	163 (91%)	17 (9%)	8	35
31	BS04	180/180 (100%)	162 (90%)	18 (10%)	7	32
32	AS05	116/116 (100%)	101 (87%)	15 (13%)	4	23
32	BS05	116/116 (100%)	106 (91%)	10 (9%)	10	40
33	AS06	90/90 (100%)	82 (91%)	8 (9%)	9	38
33	BS06	90/90 (100%)	83 (92%)	7 (8%)	12	42
34	AS07	126/126 (100%)	121 (96%)	5 (4%)	31	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	BS07	126/126 (100%)	121 (96%)	5 (4%)	31	59
35	AS08	119/119 (100%)	108 (91%)	11 (9%)	9	35
35	BS08	119/119 (100%)	112 (94%)	7 (6%)	19	51
36	AS09	98/98 (100%)	91 (93%)	7 (7%)	14	45
36	BS09	98/98 (100%)	91 (93%)	7 (7%)	14	45
37	AS10	88/88 (100%)	79 (90%)	9 (10%)	7	31
37	BS10	88/88 (100%)	80 (91%)	8 (9%)	9	36
38	AS11	86/86 (100%)	79 (92%)	7 (8%)	11	41
38	BS11	86/86 (100%)	77 (90%)	9 (10%)	7	30
39	AS12	103/103 (100%)	92 (89%)	11 (11%)	6	30
39	BS12	103/103 (100%)	93 (90%)	10 (10%)	8	33
40	AS13	94/94 (100%)	92 (98%)	2 (2%)	53	74
40	BS13	94/94 (100%)	90 (96%)	4 (4%)	29	58
41	AS14	49/49 (100%)	45 (92%)	4 (8%)	11	40
41	BS14	49/49 (100%)	46 (94%)	3 (6%)	18	50
42	AS15	79/79 (100%)	71 (90%)	8 (10%)	7	32
42	BS15	79/79 (100%)	68 (86%)	11 (14%)	3	21
43	AS16	72/72 (100%)	63 (88%)	9 (12%)	4	24
43	BS16	72/72 (100%)	68 (94%)	4 (6%)	21	52
44	AS17	94/94 (100%)	85 (90%)	9 (10%)	8	34
44	BS17	94/94 (100%)	87 (93%)	7 (7%)	13	44
45	AS18	61/61 (100%)	58 (95%)	3 (5%)	25	55
45	BS18	61/61 (100%)	53 (87%)	8 (13%)	4	23
46	AS19	69/69 (100%)	63 (91%)	6 (9%)	10	38
46	BS19	69/69 (100%)	62 (90%)	7 (10%)	7	32
47	AS20	76/76 (100%)	72 (95%)	4 (5%)	22	54
47	BS20	76/76 (100%)	72 (95%)	4 (5%)	22	54
48	ATHX	19/19 (100%)	18 (95%)	1 (5%)	22	54
48	BTHX	19/19 (100%)	19 (100%)	0	100	100
49	AL31	27/27 (100%)	26 (96%)	1 (4%)	34	62
49	BL31	27/27 (100%)	25 (93%)	2 (7%)	13	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	9618/9618 (100%)	8530 (89%)	1088 (11%)	6 28

5 of 1088 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
42	AS15	41	GLU
3	BL04	95	ARG
36	BS09	95	LYS
44	AS17	7	THR
1	BL02	150	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
50	A16S	1503/1506 (99%)	247 (16%)	21 (1%)
50	B16S	1503/1506 (99%)	249 (16%)	21 (1%)
51	A23S	2872/2879 (99%)	541 (18%)	29 (1%)
51	B23S	2872/2879 (99%)	552 (19%)	26 (0%)
52	A5S	118/119 (99%)	18 (15%)	0
52	B5S	118/119 (99%)	15 (12%)	0
53	AIRE	30/196 (15%)	4 (13%)	0
53	BIRE	30/196 (15%)	4 (13%)	0
All	All	9046/9400 (96%)	1630 (18%)	97 (1%)

5 of 1630 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
50	A16S	6	G
50	A16S	7	G
50	A16S	9	G
50	A16S	16	A
50	A16S	17	U

5 of 97 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
51	A23S	2097	C

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Mol	Chain	Res	Type
50	B16S	428	G
51	B23S	1558	A
51	A23S	2428	G
51	A23S	2719	G

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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	AL02	271/271 (100%)	0.09	4 (1%) 73 66	54, 99, 150, 228	0
1	BL02	271/271 (100%)	-0.01	1 (0%) 92 89	42, 88, 128, 197	0
2	AL03	204/204 (100%)	0.39	7 (3%) 45 37	63, 121, 179, 294	0
2	BL03	204/204 (100%)	0.16	6 (2%) 51 42	52, 114, 168, 253	0
3	AL04	202/202 (100%)	0.04	4 (1%) 65 58	60, 113, 184, 265	0
3	BL04	202/202 (100%)	0.03	3 (1%) 73 66	52, 109, 174, 242	0
4	AL05	181/181 (100%)	1.30	54 (29%) 0 0	134, 212, 272, 318	0
4	BL05	181/181 (100%)	1.42	51 (28%) 0 0	149, 227, 281, 309	0
5	AL06	159/159 (100%)	1.68	60 (37%) 0 0	124, 217, 285, 314	0
5	BL06	159/159 (100%)	0.40	9 (5%) 23 19	91, 142, 199, 240	0
6	AL09	145/145 (100%)	1.17	44 (30%) 0 0	82, 185, 244, 324	0
6	BL09	145/145 (100%)	0.75	21 (14%) 2 2	76, 156, 202, 244	0
7	AL11	147/147 (100%)	3.85	118 (80%) 0 0	256, 322, 358, 387	0
7	BL11	147/147 (100%)	4.54	113 (76%) 0 0	225, 321, 363, 388	0
8	AL13	137/137 (100%)	0.32	8 (5%) 23 18	82, 133, 194, 207	0
8	BL13	137/137 (100%)	0.34	6 (4%) 34 29	69, 120, 186, 258	0
9	AL14	122/122 (100%)	0.08	0 100 100	63, 111, 148, 180	0
9	BL14	122/122 (100%)	0.34	3 (2%) 57 49	65, 101, 143, 171	0
10	AL15	146/146 (100%)	0.45	8 (5%) 25 21	63, 138, 225, 281	0
10	BL15	146/146 (100%)	0.55	13 (8%) 9 8	46, 134, 208, 283	0
11	AL16	134/134 (100%)	0.46	10 (7%) 14 11	85, 136, 201, 289	0
11	BL16	134/134 (100%)	0.55	11 (8%) 11 9	77, 126, 219, 300	0
12	AL17	117/117 (100%)	0.52	8 (6%) 17 13	64, 113, 171, 197	0
12	BL17	117/117 (100%)	0.49	2 (1%) 70 62	67, 105, 162, 218	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AL18	98/98 (100%)	1.15	26 (26%) 0 0	125, 187, 243, 293	0
13	BL18	98/98 (100%)	0.72	17 (17%) 1 1	140, 201, 253, 275	0
14	AL19	137/137 (100%)	0.27	7 (5%) 28 24	78, 132, 216, 266	0
14	BL19	137/137 (100%)	0.15	4 (2%) 51 42	74, 124, 226, 284	0
15	AL20	117/117 (100%)	0.04	2 (1%) 70 62	66, 113, 181, 243	0
15	BL20	117/117 (100%)	-0.17	0 100 100	62, 112, 172, 232	0
16	AL21	101/101 (100%)	0.27	2 (1%) 65 58	86, 140, 206, 307	0
16	BL21	101/101 (100%)	0.18	5 (4%) 28 25	75, 130, 202, 283	0
17	AL22	112/112 (100%)	0.60	4 (3%) 42 35	71, 104, 172, 250	0
17	BL22	112/112 (100%)	0.30	3 (2%) 54 45	57, 95, 152, 210	0
18	AL23	92/92 (100%)	0.13	0 100 100	77, 117, 171, 217	0
18	BL23	92/92 (100%)	0.37	2 (2%) 62 54	62, 97, 139, 182	0
19	AL24	100/100 (100%)	1.14	20 (20%) 1 1	103, 144, 262, 296	0
19	BL24	100/100 (100%)	0.82	10 (10%) 7 6	83, 129, 234, 266	0
20	AL25	187/187 (100%)	0.96	41 (21%) 0 1	119, 190, 253, 309	0
20	BL25	187/187 (100%)	0.35	14 (7%) 14 11	111, 179, 225, 276	0
21	AL27	76/76 (100%)	0.34	5 (6%) 18 14	91, 135, 187, 233	0
21	BL27	76/76 (100%)	0.42	5 (6%) 18 14	88, 132, 198, 249	0
22	AL28	88/88 (100%)	0.35	3 (3%) 45 37	61, 109, 186, 283	0
22	BL28	88/88 (100%)	0.37	4 (4%) 33 28	61, 105, 182, 248	0
23	AL29	62/62 (100%)	0.33	4 (6%) 18 14	99, 137, 237, 255	0
23	BL29	62/62 (100%)	0.28	6 (9%) 7 7	59, 108, 218, 253	0
24	AL30	59/59 (100%)	1.40	15 (25%) 0 0	101, 132, 214, 248	0
24	BL30	59/59 (100%)	0.92	5 (8%) 10 9	78, 131, 207, 286	0
25	AL32	52/52 (100%)	0.51	7 (13%) 3 3	66, 117, 205, 239	0
25	BL32	52/52 (100%)	0.06	2 (3%) 40 33	54, 105, 215, 269	0
26	AL33	44/44 (100%)	9.55	44 (100%) 0 0	140, 241, 282, 301	0
26	BL33	44/44 (100%)	5.13	39 (88%) 0 0	193, 248, 287, 304	0
27	AL34	48/48 (100%)	-0.04	0 100 100	59, 89, 129, 203	0
27	BL34	48/48 (100%)	0.14	0 100 100	41, 71, 113, 175	0
28	AL35	63/63 (100%)	0.21	1 (1%) 72 64	69, 119, 192, 227	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	BL35	63/63 (100%)	0.16	1 (1%) 72 64	71, 114, 188, 225	0
29	AS02	234/234 (100%)	0.56	22 (9%) 8 7	129, 202, 274, 300	0
29	BS02	234/234 (100%)	0.64	26 (11%) 5 5	113, 213, 279, 309	0
30	AS03	206/206 (100%)	0.44	15 (7%) 15 12	124, 200, 264, 308	0
30	BS03	206/206 (100%)	0.34	12 (5%) 23 18	119, 196, 253, 298	0
31	AS04	208/208 (100%)	0.23	5 (2%) 59 50	99, 152, 216, 239	0
31	BS04	208/208 (100%)	0.73	31 (14%) 2 2	128, 204, 275, 307	0
32	AS05	151/151 (100%)	0.11	4 (2%) 56 47	94, 144, 199, 262	0
32	BS05	151/151 (100%)	0.04	5 (3%) 46 38	105, 164, 216, 261	0
33	AS06	101/101 (100%)	0.17	5 (4%) 28 25	122, 172, 234, 258	0
33	BS06	101/101 (100%)	-0.10	2 (1%) 65 58	96, 149, 203, 230	0
34	AS07	155/155 (100%)	0.91	32 (20%) 1 1	154, 214, 272, 300	0
34	BS07	155/155 (100%)	0.95	30 (19%) 1 1	151, 222, 282, 309	0
35	AS08	138/138 (100%)	0.47	12 (8%) 10 8	92, 142, 195, 229	0
35	BS08	138/138 (100%)	0.58	14 (10%) 7 6	95, 168, 215, 256	0
36	AS09	127/127 (100%)	1.61	38 (29%) 0 0	139, 274, 335, 355	0
36	BS09	127/127 (100%)	1.31	31 (24%) 0 0	150, 262, 310, 343	0
37	AS10	98/98 (100%)	2.25	49 (50%) 0 0	138, 260, 344, 356	0
37	BS10	98/98 (100%)	1.61	33 (33%) 0 0	142, 246, 308, 339	0
38	AS11	114/114 (100%)	0.43	13 (11%) 5 5	95, 150, 210, 267	0
38	BS11	114/114 (100%)	0.11	3 (2%) 56 47	85, 138, 193, 251	0
39	AS12	122/122 (100%)	0.72	12 (9%) 7 7	74, 121, 166, 201	0
39	BS12	122/122 (100%)	0.53	8 (6%) 18 14	105, 147, 192, 230	0
40	AS13	117/117 (100%)	1.65	41 (35%) 0 0	152, 259, 315, 352	0
40	BS13	117/117 (100%)	1.11	30 (25%) 0 0	159, 253, 318, 340	0
41	AS14	60/60 (100%)	1.05	11 (18%) 1 1	148, 194, 241, 329	0
41	BS14	60/60 (100%)	1.02	9 (15%) 2 2	144, 196, 243, 301	0
42	AS15	88/88 (100%)	0.32	5 (5%) 23 19	100, 139, 183, 229	0
42	BS15	88/88 (100%)	0.48	5 (5%) 23 19	92, 146, 201, 215	0
43	AS16	83/83 (100%)	0.84	12 (14%) 2 2	104, 137, 187, 233	0
43	BS16	83/83 (100%)	1.44	27 (32%) 0 0	148, 199, 260, 310	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	AS17	99/99 (100%)	0.15	3 (3%) 50 40	92, 125, 166, 223	0
44	BS17	99/99 (100%)	0.36	3 (3%) 50 40	105, 149, 191, 235	0
45	AS18	70/70 (100%)	0.85	9 (12%) 3 4	110, 174, 234, 253	0
45	BS18	70/70 (100%)	0.49	1 (1%) 75 68	100, 150, 195, 247	0
46	AS19	78/78 (100%)	2.27	38 (48%) 0 0	170, 255, 301, 353	0
46	BS19	78/78 (100%)	1.70	28 (35%) 0 0	172, 257, 302, 314	0
47	AS20	99/99 (100%)	0.69	9 (9%) 9 7	106, 149, 228, 256	0
47	BS20	99/99 (100%)	1.18	22 (22%) 0 0	120, 178, 249, 291	0
48	ATHX	24/24 (100%)	0.68	1 (4%) 36 30	153, 263, 279, 301	0
48	BTHX	24/24 (100%)	1.20	8 (33%) 0 0	209, 276, 325, 350	0
49	AL31	30/30 (100%)	1.85	13 (43%) 0 0	206, 272, 323, 351	0
49	BL31	30/30 (100%)	1.81	11 (36%) 0 0	216, 303, 332, 352	0
50	A16S	1504/1506 (99%)	0.19	50 (3%) 46 38	76, 144, 270, 413	0
50	B16S	1504/1506 (99%)	0.23	56 (3%) 41 34	79, 171, 296, 397	0
51	A23S	2876/2879 (99%)	0.23	126 (4%) 34 29	58, 115, 252, 440	0
51	B23S	2876/2879 (99%)	0.14	117 (4%) 37 31	52, 103, 243, 414	0
52	A5S	119/119 (100%)	-0.03	1 (0%) 86 81	122, 178, 230, 282	0
52	B5S	119/119 (100%)	-0.19	1 (0%) 86 81	124, 178, 236, 309	0
53	AIRE	32/196 (16%)	1.84	10 (31%) 0 0	131, 266, 505, 516	0
53	BIRE	32/196 (16%)	0.48	2 (6%) 20 15	145, 222, 413, 467	0
All	All	20668/21006 (98%)	0.49	1913 (9%) 8 7	41, 141, 282, 516	0

The worst 5 of 1913 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
7	BL11	1	MET	31.2
26	AL33	49	HIS	20.9
26	AL33	13	CYS	20.3
26	AL33	47	THR	17.4
26	AL33	22	ALA	16.8

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

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

6.3 Carbohydrates [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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
54	ZN	BS04	301	1/1	0.90	0.23	100,100,100,100	0
54	ZN	AS14	101	1/1	0.95	0.13	120,120,120,120	0
54	ZN	BS14	101	1/1	0.96	0.10	137,137,137,137	0
54	ZN	AS04	301	1/1	0.99	0.24	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.