



wwPDB EM Validation Summary Report ⓘ

Dec 11, 2022 – 05:28 pm GMT

PDB ID : 5AJ0
EMDB ID : EMD-2875
Title : Cryo electron microscopy of actively translating human polysomes (POST state).
Authors : Behrmann, E.; Loerke, J.; Budkevich, T.V.; Yamamoto, K.; Schmidt, A.; Penczek, P.A.; Vos, M.R.; Burger, J.; Mielke, T.; Scheerer, P.; Spahn, C.M.T.
Deposited on : 2015-02-19
Resolution : 3.50 Å (reported)
Based on initial model : 4UJE

This is a wwPDB EM 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/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

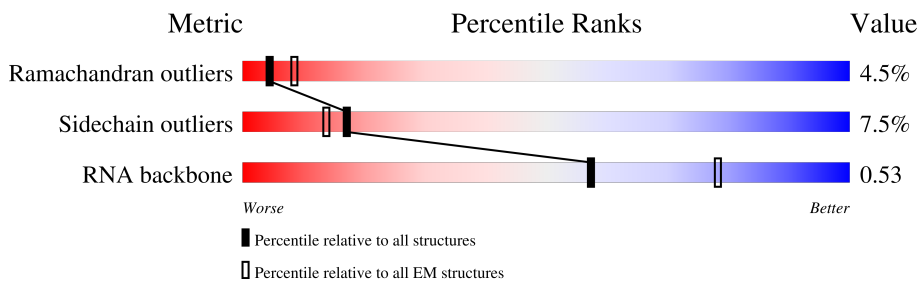
EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A3	194	81% (Upper red bar) 46% (Green), 27% (Yellow), 8% (Orange), 19% (Grey)
2	A4	121	98% (Upper red bar) 44% (Green), 49% (Yellow), 6% (Orange), 1% (Grey)
3	AA	257	98% (Upper red bar) 92% (Green), 6% (Yellow), 2% (Orange), 1% (Grey)
4	AB	403	98% (Upper red bar) 90% (Green), 7% (Yellow), 3% (Orange), 1% (Grey)
5	AC	427	85% (Upper red bar) 76% (Green), 8% (Yellow), 15% (Grey)
6	AD	297	99% (Upper red bar) 91% (Green), 8% (Yellow), 1% (Orange), 1% (Grey)
7	AE	288	67% (Upper red bar) 46% (Green), 18% (Yellow), 33% (Grey)
8	AF	248	94% (Upper red bar) 88% (Green), 6% (Yellow), 6% (Grey)

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Mol	Chain	Length	Quality of chain
9	AG	266	88% 84% 12%
10	AH	192	99% 91% 7%
11	AI	214	97% 93% 5%
12	AJ	178	95% 90% 5%
13	AK	317	34% 17% 11% 5% 66%
14	AL	211	97% 82% 12%
15	AM	215	65% 56% 8% 35%
16	AN	204	100% 95% 5%
17	AO	203	96% 93%
18	AP	184	83% 77% 6% 17%
19	AQ	188	99% 81% 16%
20	AR	196	92% 83% 8% 8%
21	AS	176	99% 92% 6%
22	AT	160	98% 88% 11%
23	AU	128	77% 62% 14% 23%
24	AV	140	92% 91% 8%
25	AW	157	77% 68% 9% 23%
26	AX	156	75% 69% 6% 25%
27	AY	145	88% 78% 10% 12%
28	AZ	136	99% 82% 14%
29	Aa	148	99% 90% 9%
30	Ab	159	43% 38% 5% 57%
31	Ac	115	90% 85% 10%
32	Ad	125	85% 82% 15%
33	Ae	135	96% 91%

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Mol	Chain	Length	Quality of chain
34	Af	110	99% 90% 6% ...
35	Ag	117	97% 91% 6% .
36	Ah	123	99% 83% 16% .
37	Ai	105	92% 85% 8% 8%
38	Aj	97	87% 72% 12% . 13%
39	Ak	70	99% 66% 24% 9% .
40	Al	51	98% 92% 6% .
41	Am	128	39% 38% . 61%
42	An	25	100% 96% .
43	Ao	106	99% 88% 11% .
44	Ap	92	99% 93% 5% .
45	Aq	165	84% 41% 30% 11% . 16%
46	At	137	89% 74% 15% 11%
47	Au	217	100% 87% 10% . .
48	A2	5029	72% 42% 26% . 28%
49	B1	1869	91% 54% 34% . 9%
50	BA	295	73% 66% 6% . 27%
51	BB	264	80% 68% 12% 20%
52	BC	293	76% 69% 7% 24%
53	BD	243	91% 84% 6% 9%
54	BE	263	98% 90% 6% . .
55	BF	204	93% 85% 7% 7%
56	BG	249	93% 80% 14% 7%
57	BH	194	94% 90% 5% 6%
58	BI	208	100% 93% 7%

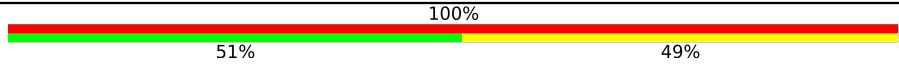
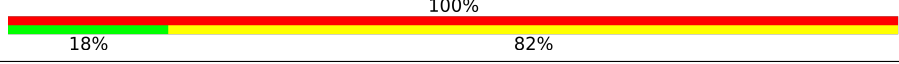
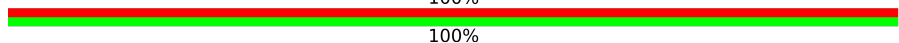
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Mol	Chain	Length	Quality of chain
59	BJ	194	92% 81% 10% 8%
60	BK	165	59% 54% 41%
61	BL	158	97% 89% 8%
62	BM	132	91% 87% 9%
63	BN	151	99% 86% 12%
64	BO	151	90% 79% 10% 10%
65	BP	145	83% 65% 12% 17%
66	BQ	146	95% 79% 14% 5%
67	BR	135	93% 74% 15% 7%
68	BS	152	91% 77% 14% 9%
69	BT	145	99% 88% 10%
70	BU	119	82% 75% 7% 18%
71	BV	83	98% 89% 8%
72	BW	130	99% 92% 5%
73	BX	143	97% 86% 10%
74	BY	133	94% 90% 6%
75	BZ	125	69% 52% 17% 31%
76	Ba	115	84% 76% 8% 16%
77	Bb	84	95% 93% 5%
78	Bc	69	90% 84% 10%
79	Bd	56	91% 71% 18% 9%
80	Be	59	93% 80% 14% 7%
81	Bf	156	47% 46% 53%
82	Bg	317	99% 96%
83	Bv	76	100% 46% 51%

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Mol	Chain	Length	Quality of chain
83	Bw	76	 100% 51% 49%
84	Bx	28	 100% 18% 82%
85	By	24	 100% 100%

2 Entry composition [i](#)

There are 87 unique types of molecules in this entry. The entry contains 218559 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 5.8S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	A3	157	3337	1489	587	1104	157	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	A4	119	2541	1132	454	836	119	0	0

- Molecule 3 is a protein called 60S ribosomal protein L8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	AA	252	1930	1209	395	320	6	0	0

- Molecule 4 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	AB	394	3178	2024	596	544	14	0	0

- Molecule 5 is a protein called 60S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	AC	363	2888	1817	577	480	14	0	0

- Molecule 6 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	AD	294	2392	1510	436	432	14	0	0

- Molecule 7 is a protein called 60S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	AE	194	1571	1013	294	263	1	0	0

- Molecule 8 is a protein called 60S ribosomal protein L7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	AF	234	1950	1252	376	313	9	0	0

- Molecule 9 is a protein called 60S ribosomal protein L7a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	AG	234	1880	1197	362	317	4	0	0

- Molecule 10 is a protein called 60S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	AH	191	1526	960	285	275	6	0	0

- Molecule 11 is a protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	AI	208	1692	1074	327	278	13	0	0

- Molecule 12 is a protein called 60S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	AJ	169	1353	855	252	240	6	0	0

- Molecule 13 is a protein called 60S acidic ribosomal protein P0.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	AK	109	872	554	159	151	8	0	0

- Molecule 14 is a protein called 60S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	AL	205	Total	C	N	O	S	0	0
			1657	1036	344	273	4		

- Molecule 15 is a protein called 60S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	AM	139	Total	C	N	O	S	0	0
			1138	730	218	183	7		

- Molecule 16 is a protein called 60S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	AN	203	Total	C	N	O	S	0	0
			1701	1072	359	266	4		

- Molecule 17 is a protein called 60S ribosomal protein L13a.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	AO	195	Total	C	N	O	S	0	0
			1606	1034	315	252	5		

- Molecule 18 is a protein called 60S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	AP	153	Total	C	N	O	S	0	0
			1242	776	241	216	9		

- Molecule 19 is a protein called 60S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	AQ	187	Total	C	N	O	S	0	0
			1513	944	314	250	5		

- Molecule 20 is a protein called 60S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	AR	181	Total	C	N	O	S	0	0
			1517	938	329	241	9		

- Molecule 21 is a protein called 60S ribosomal protein L18a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
21	AS	175	1449	921	283	234	11	0	0

- Molecule 22 is a protein called 60S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
22	AT	157	1284	815	250	214	5	0	0

- Molecule 23 is a protein called 60S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
23	AU	99	808	518	141	147	2	0	0

- Molecule 24 is a protein called 60S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
24	AV	129	969	613	182	169	5	0	0

- Molecule 25 is a protein called 60S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
25	AW	121	989	617	202	167	3	0	0

- Molecule 26 is a protein called 60S ribosomal protein L23a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
26	AX	117	958	612	180	165	1	0	0

- Molecule 27 is a protein called 60S ribosomal protein L26.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
27	AY	127	1064	668	216	177	3	0	0

- Molecule 28 is a protein called 60S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
28	AZ	134	1103	712	207	181	3	0	0

- Molecule 29 is a protein called 60S ribosomal protein L27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
29	Aa	147	1162	736	237	186	3	0	0

- Molecule 30 is a protein called 60S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
30	Ab	68	559	344	122	90	3	0	0

- Molecule 31 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
31	Ac	103	801	508	141	145	7	0	0

- Molecule 32 is a protein called 60S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
32	Ad	106	879	555	170	152	2	0	0

- Molecule 33 is a protein called 60S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
33	Ae	129	1064	673	220	166	5	0	0

- Molecule 34 is a protein called 60S ribosomal protein L35a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
34	Af	109	876	555	174	144	3	0	0

- Molecule 35 is a protein called 60S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
35	Ag	114	906	566	187	147	6	0	0

- Molecule 36 is a protein called 60S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
36	Ah	122	1015	641	205	168	1	0	0

- Molecule 37 is a protein called 60S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
37	Ai	97	794	497	168	124	5	0	0

- Molecule 38 is a protein called 60S ribosomal protein L37.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
38	Aj	84	689	423	152	109	5	0	0

- Molecule 39 is a protein called 60S ribosomal protein L38.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
39	Ak	69	569	366	103	99	1	0	0

- Molecule 40 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
40	Al	50	444	281	98	64	1	0	0

- Molecule 41 is a protein called Ubiquitin-60S ribosomal protein L40.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
41	Am	50	411	254	87	64	6	0	0

- Molecule 42 is a protein called 60S ribosomal protein L41.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	An	25	Total	C	N	O	S	0	0
			240	145	64	28	3		

- Molecule 43 is a protein called 60S ribosomal protein L36a.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	Ao	105	Total	C	N	O	S	0	0
			863	542	175	140	6		

- Molecule 44 is a protein called 60S ribosomal protein L37a.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	Ap	91	Total	C	N	O	S	0	0
			708	445	136	120	7		

- Molecule 45 is a protein called 60S ribosomal protein L12.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	Aq	138	Total	C	N	O	S	0	0
			1046	654	196	193	3		

- Molecule 46 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	At	122	Total	C	N	O	S	0	0
			980	607	204	165	4		

- Molecule 47 is a protein called 60S ribosomal protein L10a.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	Au	217	Total	C	N	O	S	0	0
			1744	1114	314	307	9		

- Molecule 48 is a RNA chain called 28S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	A2	3612	Total	C	N	O	P	0	0
			77427	34482	14158	25175	3612		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A2	245	C	-	insertion	GB 337381
A2	246	C	-	insertion	GB 337381
A2	247	C	-	insertion	GB 337381
A2	4684	G	-	insertion	GB 337381

- Molecule 49 is a RNA chain called 18S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
49	B1	1708	36456	16274	6546	11928	1708	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
50	BA	215	1704	1083	298	315	8	0	0

- Molecule 51 is a protein called 40S ribosomal protein S3a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
51	BB	212	1722	1093	308	307	14	0	0

- Molecule 52 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
52	BC	222	1724	1114	296	304	10	0	0

- Molecule 53 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
53	BD	220	1709	1090	308	304	7	0	0

- Molecule 54 is a protein called 40S ribosomal protein S4, Y isoform 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
54	BE	257	2031	1298	381	344	8	0	0

- Molecule 55 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	BF	190	Total	C	N	O	S	0	0
			1502	939	285	271	7		

- Molecule 56 is a protein called 40S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	BG	232	Total	C	N	O	S	0	0
			1884	1176	379	322	7		

- Molecule 57 is a protein called 40S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	BH	183	Total	C	N	O	S	0	0
			1479	941	272	265	1		

- Molecule 58 is a protein called 40S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	BI	207	Total	C	N	O	S	0	0
			1696	1064	334	293	5		

- Molecule 59 is a protein called 40S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	BJ	179	Total	C	N	O	S	0	0
			1495	953	299	241	2		

- Molecule 60 is a protein called 40S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	BK	98	Total	C	N	O	S	0	0
			827	539	148	134	6		

- Molecule 61 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	BL	153	Total	C	N	O	S	0	0
			1258	804	235	213	6		

- Molecule 62 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
62	BM	120	931	584	164	174	9	0	0

- Molecule 63 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
63	BN	149	1202	770	228	203	1	0	0

- Molecule 64 is a protein called 40S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
64	BO	136	1016	621	199	190	6	0	0

- Molecule 65 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
65	BP	120	999	636	188	168	7	0	0

- Molecule 66 is a protein called 40S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
66	BQ	139	1109	704	210	192	3	0	0

- Molecule 67 is a protein called 40S ribosomal protein S17-like.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
67	BR	125	1011	634	187	186	4	0	0

- Molecule 68 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
68	BS	139	1154	725	233	195	1	0	0

- Molecule 69 is a protein called 40S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
69	BT	143	1112	697	214	198	3	0	0

- Molecule 70 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
70	BU	97	769	483	144	138	4	0	0

- Molecule 71 is a protein called 40S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
71	BV	81	617	380	114	118	5	0	0

- Molecule 72 is a protein called 40S ribosomal protein S15a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
72	BW	129	1034	659	193	176	6	0	0

- Molecule 73 is a protein called 40S ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
73	BX	139	1080	682	214	181	3	0	0

- Molecule 74 is a protein called 40S ribosomal protein S24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
74	BY	125	1015	642	199	169	5	0	0

- Molecule 75 is a protein called 40S ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
75	BZ	86	688	442	129	116	1	0	0

- Molecule 76 is a protein called 40S ribosomal protein S26.

Mol	Chain	Residues	Atoms					AltConf	Trace
76	Ba	97	Total	C	N	O	S	0	0
			774	481	160	128	5		

- Molecule 77 is a protein called 40S ribosomal protein S27.

Mol	Chain	Residues	Atoms					AltConf	Trace
77	Bb	80	Total	C	N	O	S	0	0
			625	391	116	111	7		

- Molecule 78 is a protein called 40S ribosomal protein S28.

Mol	Chain	Residues	Atoms					AltConf	Trace
78	Bc	62	Total	C	N	O	S	0	0
			488	297	97	92	2		

- Molecule 79 is a protein called 40S ribosomal protein S29.

Mol	Chain	Residues	Atoms					AltConf	Trace
79	Bd	51	Total	C	N	O	S	0	0
			427	269	87	66	5		

- Molecule 80 is a protein called 40S ribosomal protein S30.

Mol	Chain	Residues	Atoms					AltConf	Trace
80	Be	55	Total	C	N	O	S	0	0
			437	272	96	68	1		

- Molecule 81 is a protein called Ubiquitin-40S ribosomal protein S27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
81	Bf	73	Total	C	N	O	S	0	0
			601	379	115	100	7		

- Molecule 82 is a protein called Guanine nucleotide-binding protein subunit beta-2-like 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
82	Bg	314	Total	C	N	O	S	0	0
			2440	1537	425	466	12		

- Molecule 83 is a RNA chain called tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
83	Bv	76	Total	C	N	O	P	0	0
			1623	723	290	534	76		
83	Bw	76	Total	C	N	O	P	0	0
			1623	723	290	534	76		

- Molecule 84 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
84	Bx	28	Total	C	N	O	P	0	0
			561	252	56	225	28		

- Molecule 85 is a protein called Nascent protein chain.

Mol	Chain	Residues	Atoms				AltConf	Trace
85	By	24	Total	C	N	O	0	0
			120	72	24	24		

- Molecule 86 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
86	A3	8	Total	Mg	0
			8	8	
86	A4	9	Total	Mg	0
			9	9	
86	AA	1	Total	Mg	0
			1	1	
86	AB	2	Total	Mg	0
			2	2	
86	AN	2	Total	Mg	0
			2	2	
86	AY	1	Total	Mg	0
			1	1	
86	Aa	3	Total	Mg	0
			3	3	
86	Ae	2	Total	Mg	0
			2	2	
86	An	1	Total	Mg	0
			1	1	
86	A2	220	Total	Mg	0
			220	220	
86	B1	72	Total	Mg	0
			72	72	

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Mol	Chain	Residues	Atoms		AltConf
86	BD	1	Total 1	Mg 1	0
86	BX	1	Total 1	Mg 1	0
86	Bv	2	Total 2	Mg 2	0
86	Bx	1	Total 1	Mg 1	0

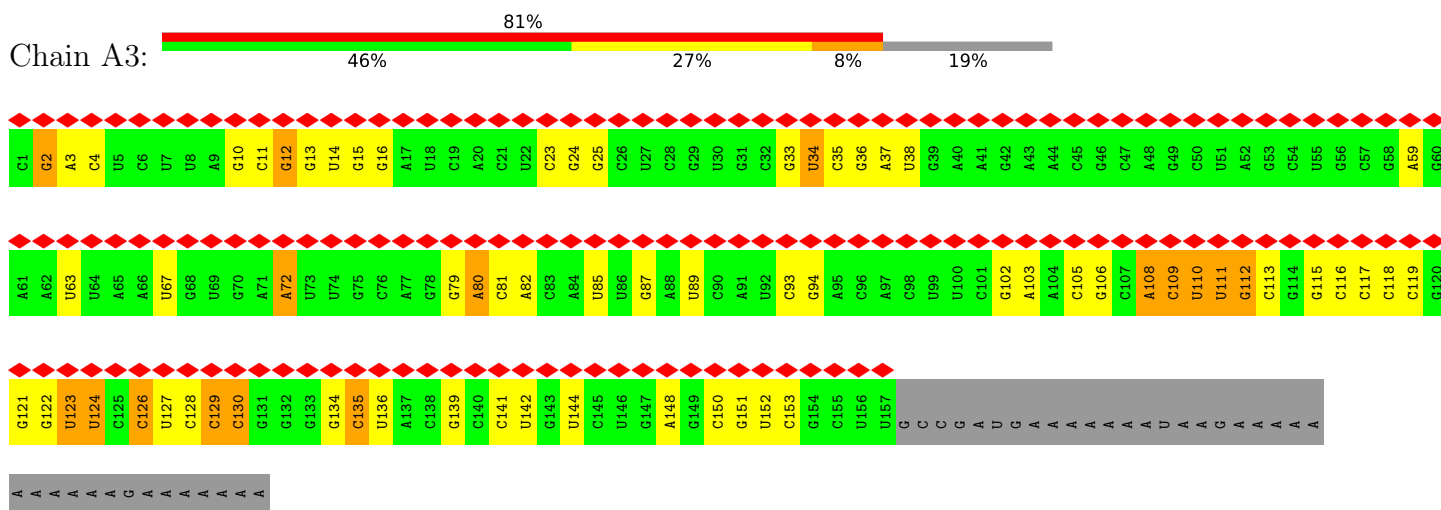
- Molecule 87 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
87	Aj	1	Total 1	Zn 1	0
87	Ao	1	Total 1	Zn 1	0
87	Ap	1	Total 1	Zn 1	0
87	Ba	1	Total 1	Zn 1	0
87	Bd	1	Total 1	Zn 1	0

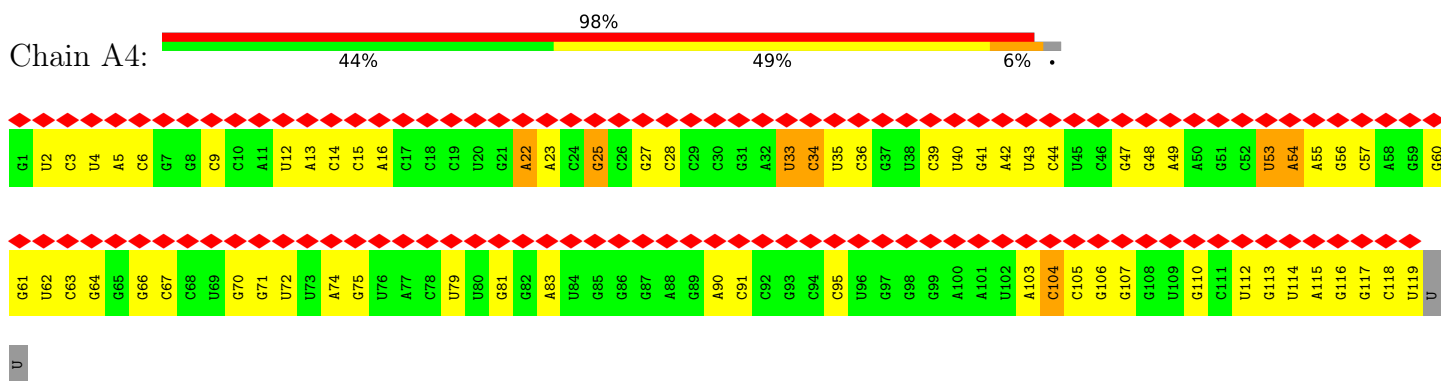
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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.

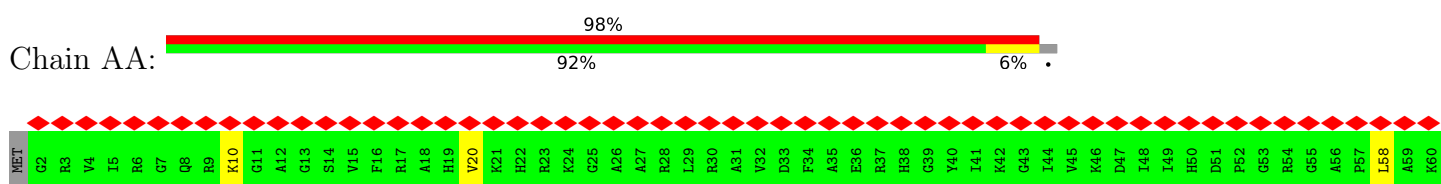
- Molecule 1: 5.8S ribosomal RNA

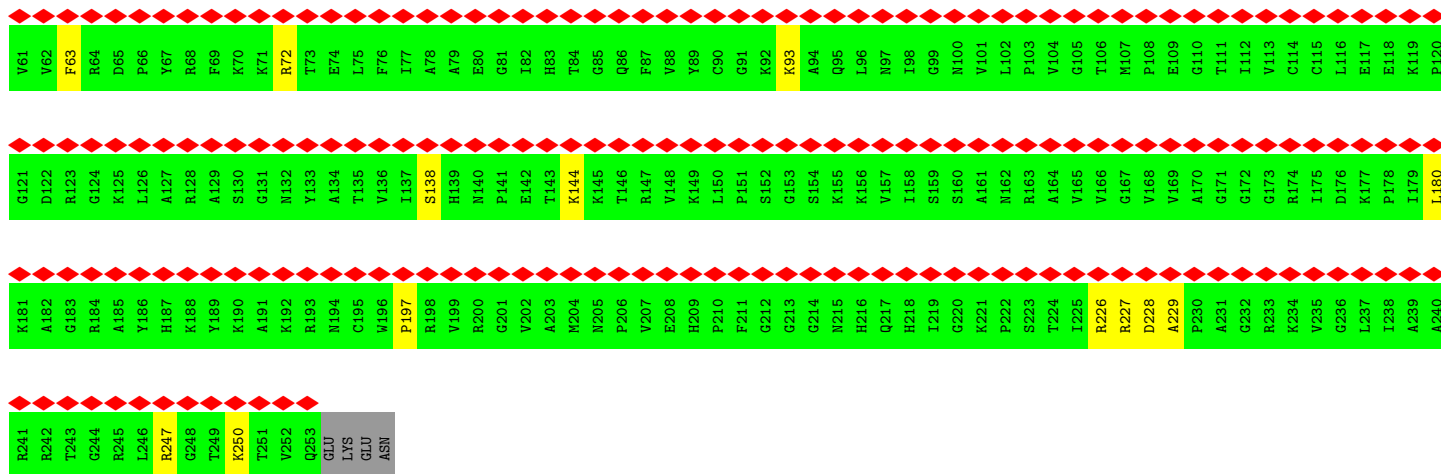


- Molecule 2: 5S ribosomal RNA

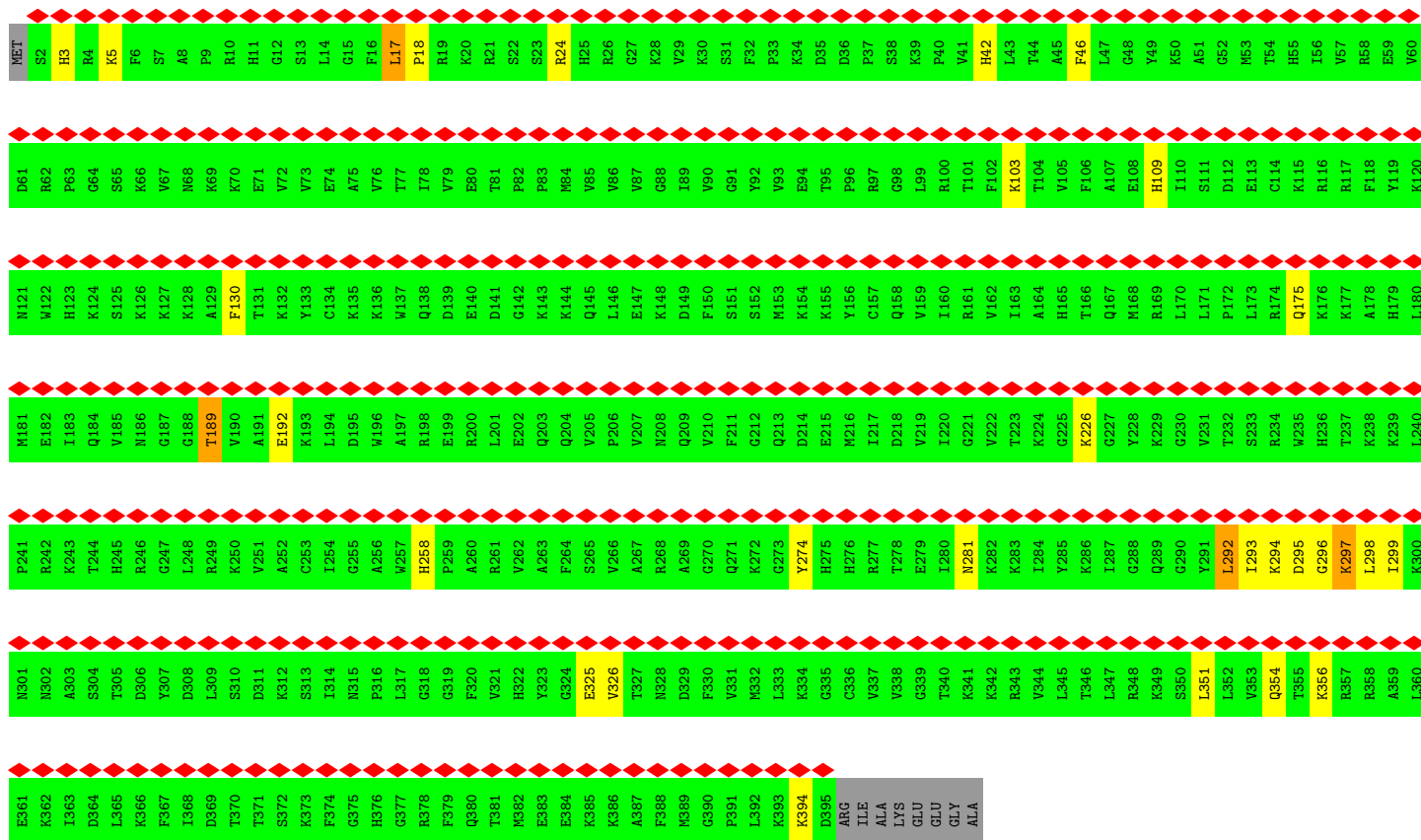
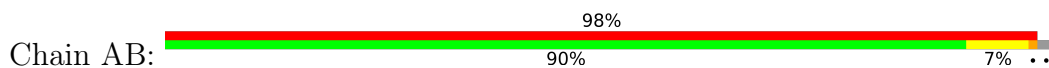


- Molecule 3: 60S ribosomal protein L8

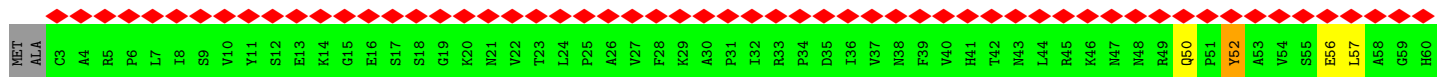
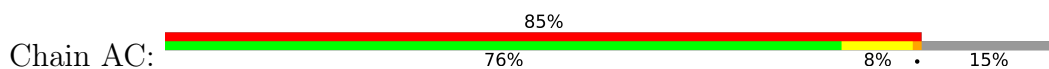


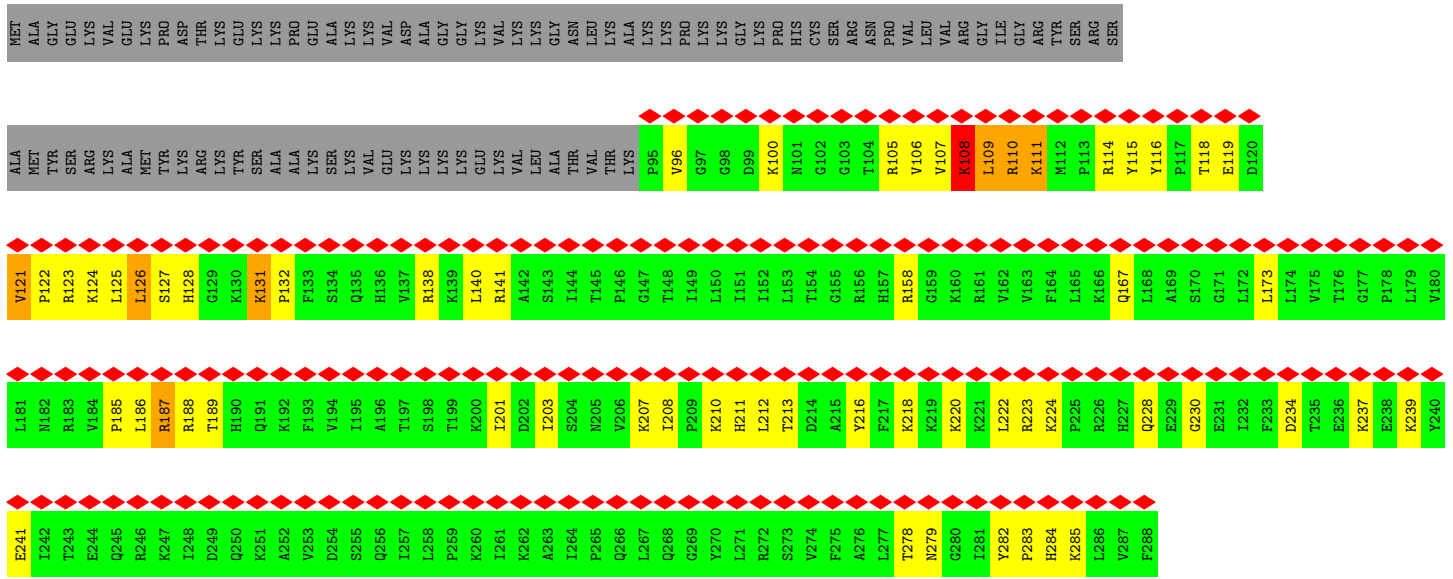


• Molecule 4: 60S ribosomal protein L3

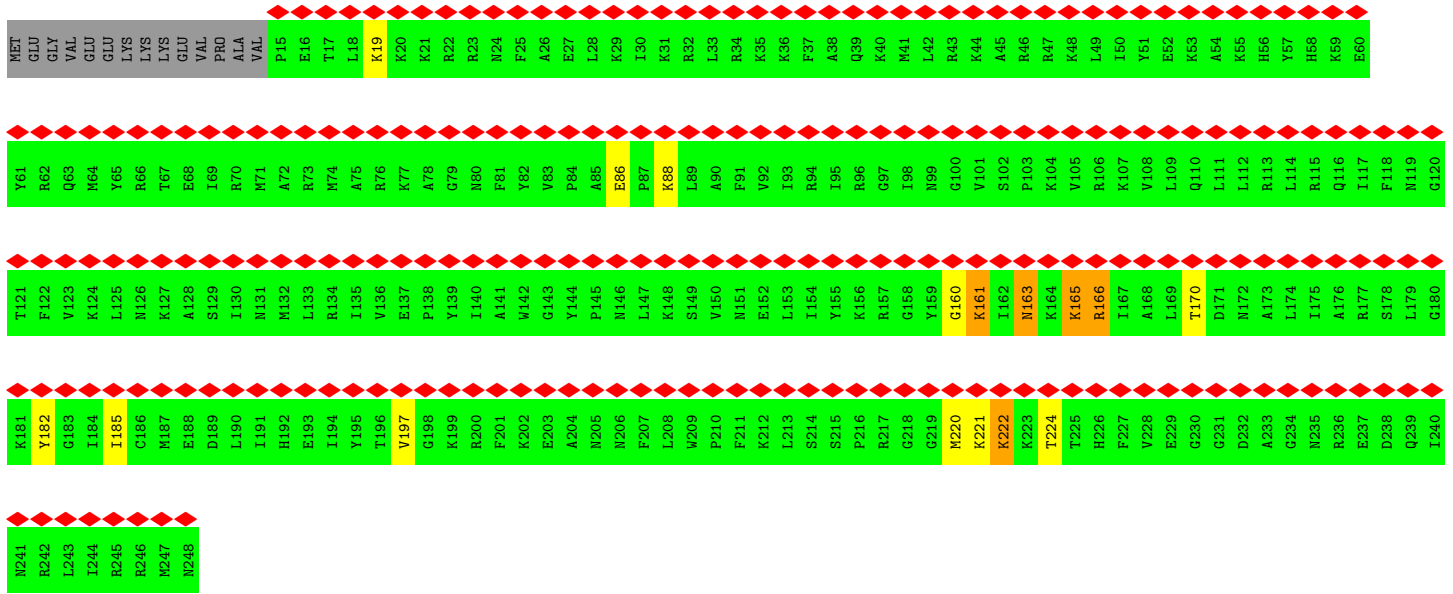
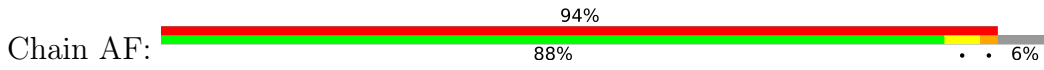


• Molecule 5: 60S ribosomal protein L4

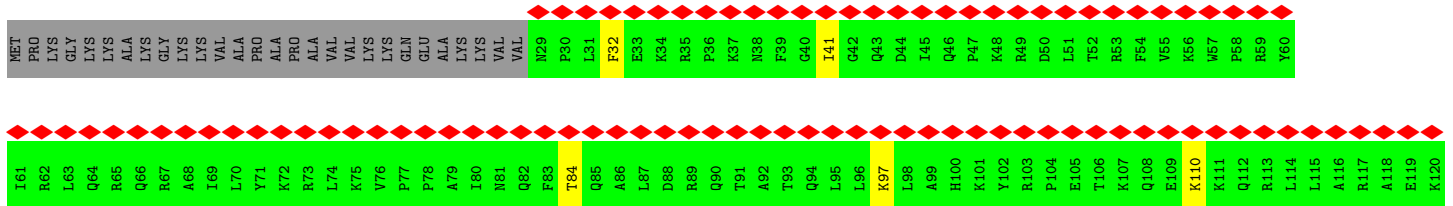
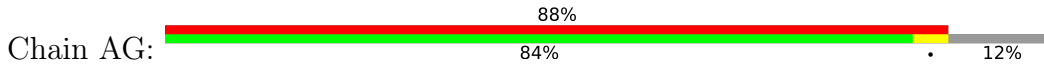


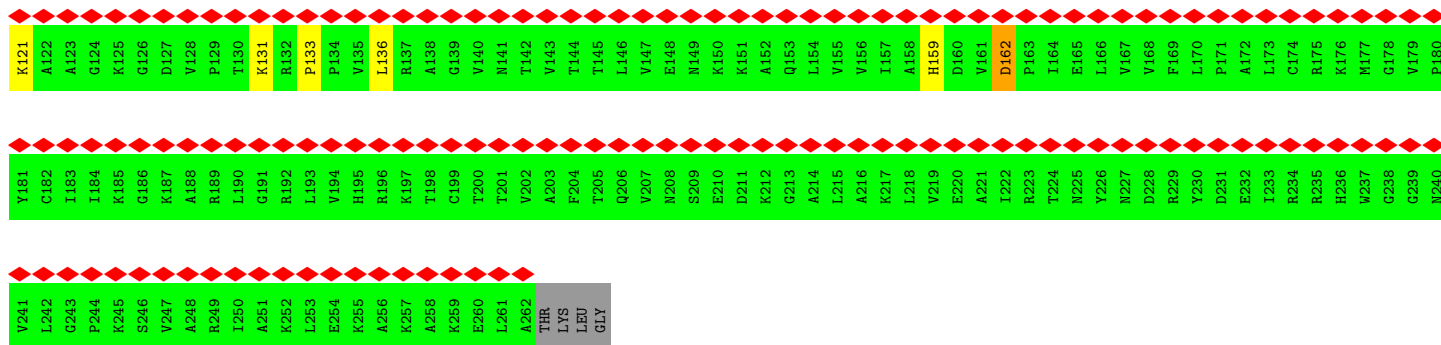


• Molecule 8: 60S ribosomal protein L7

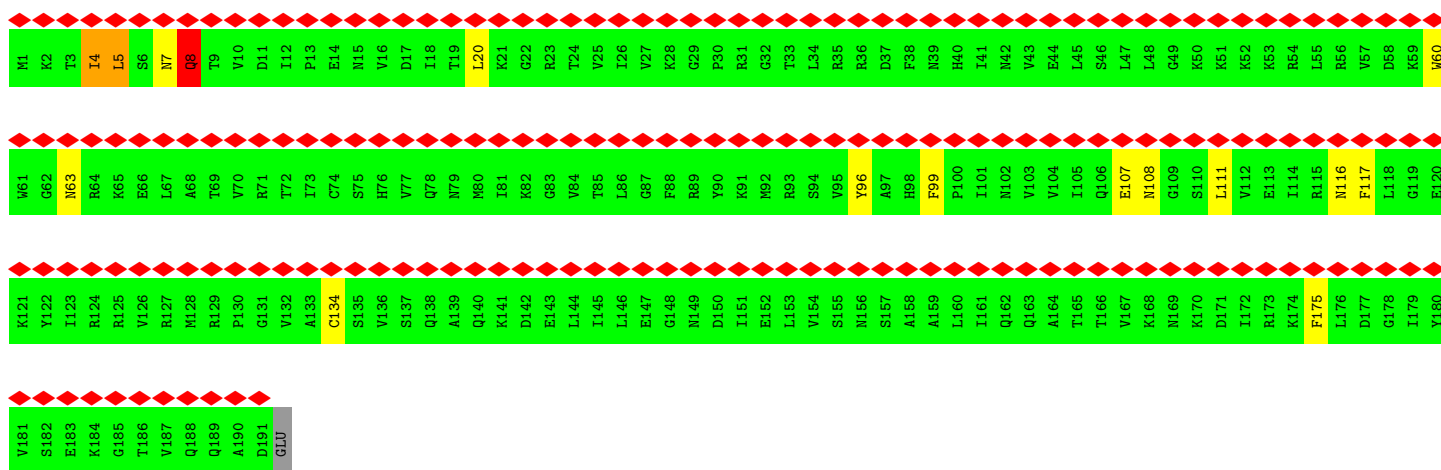
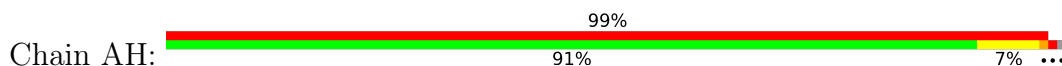


• Molecule 9: 60S ribosomal protein L7a

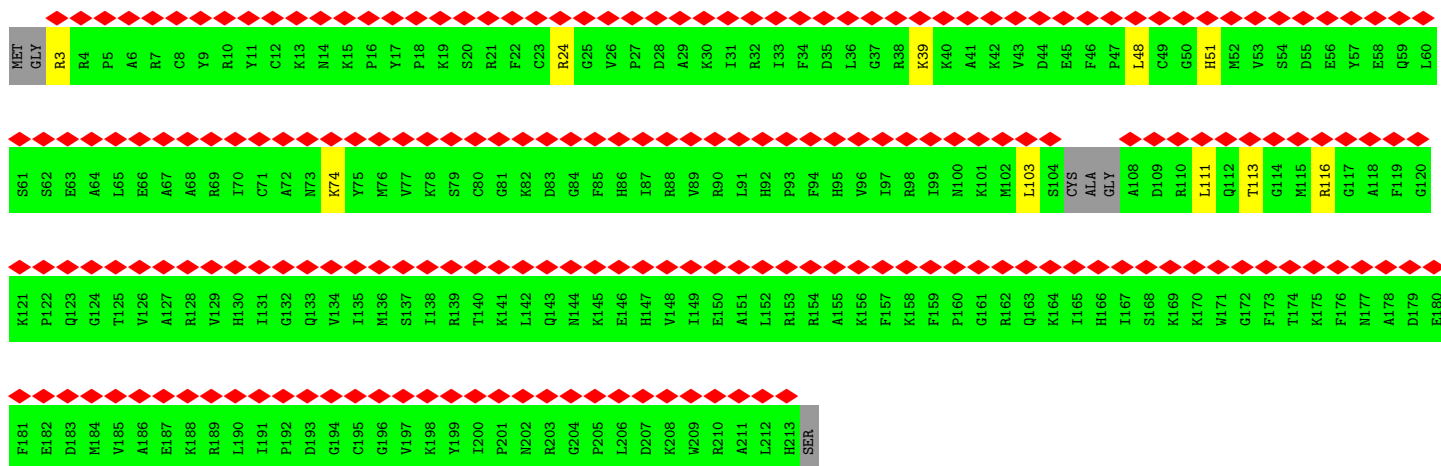
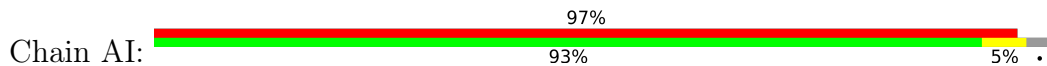




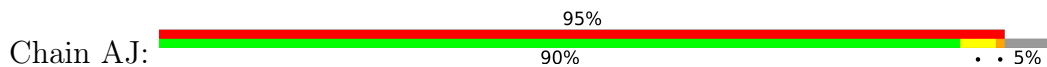
• Molecule 10: 60S ribosomal protein L9

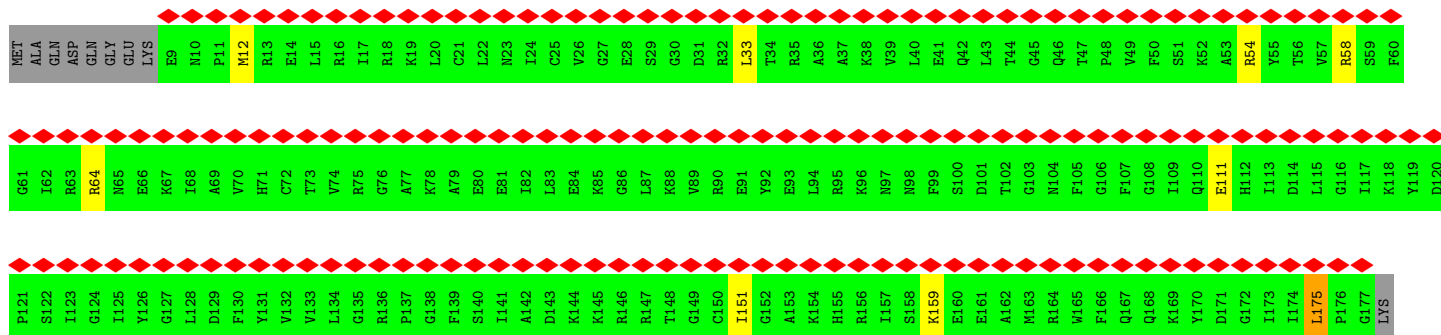


• Molecule 11: 60S ribosomal protein L10

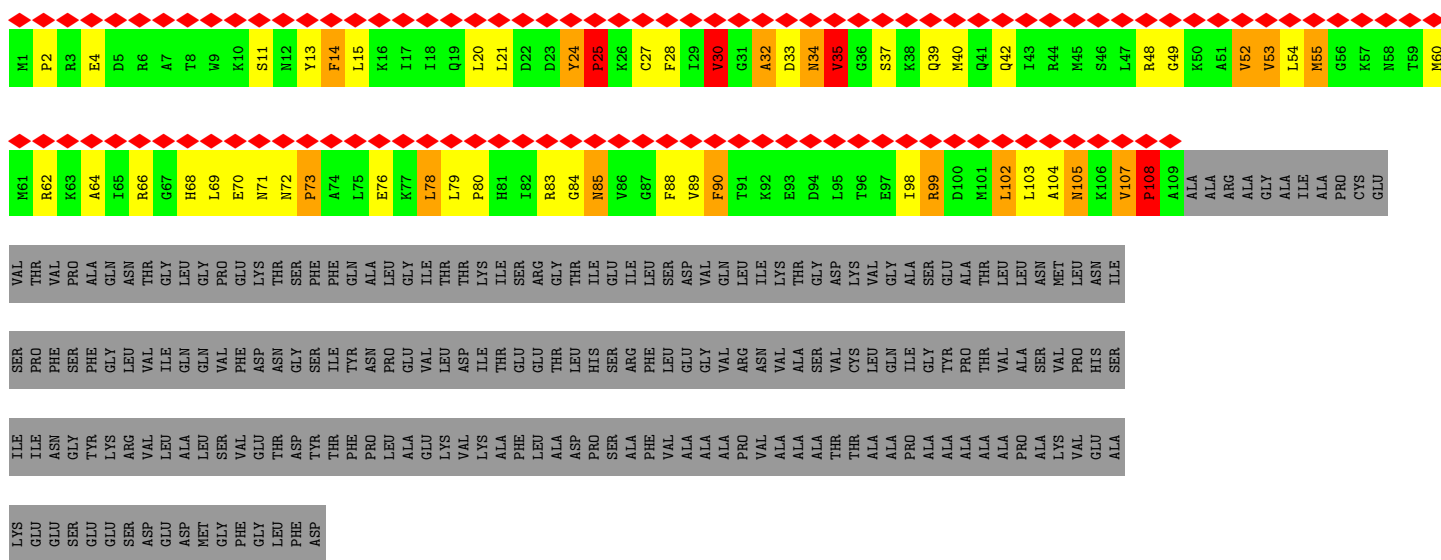


• Molecule 12: 60S ribosomal protein L11

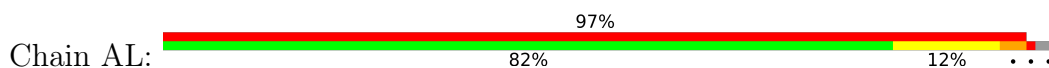




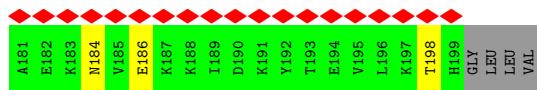
• Molecule 13: 60S acidic ribosomal protein P0



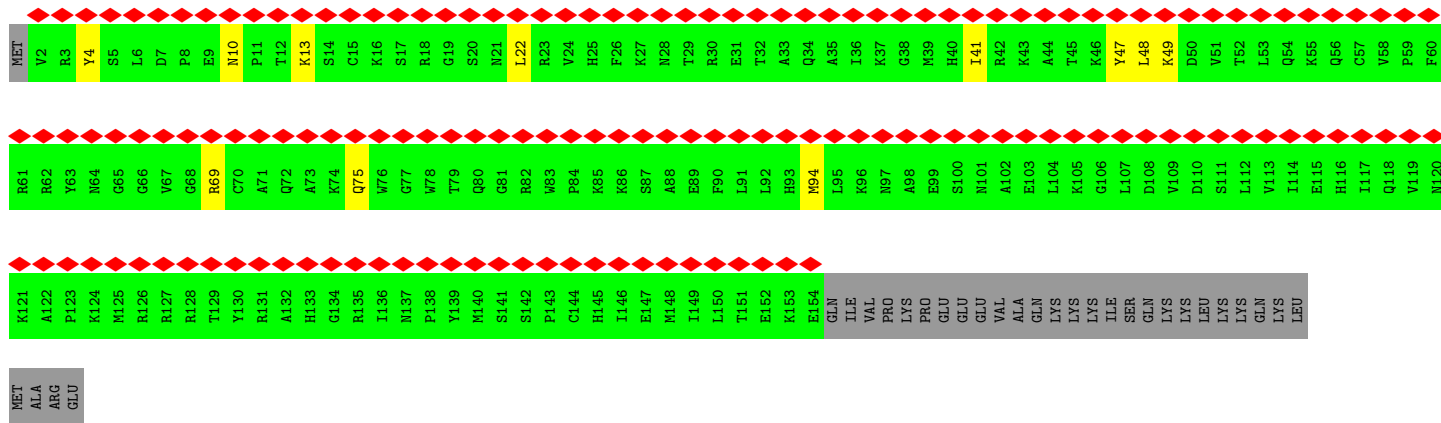
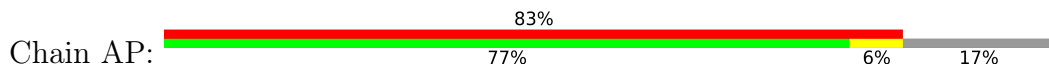
• Molecule 14: 60S ribosomal protein L13



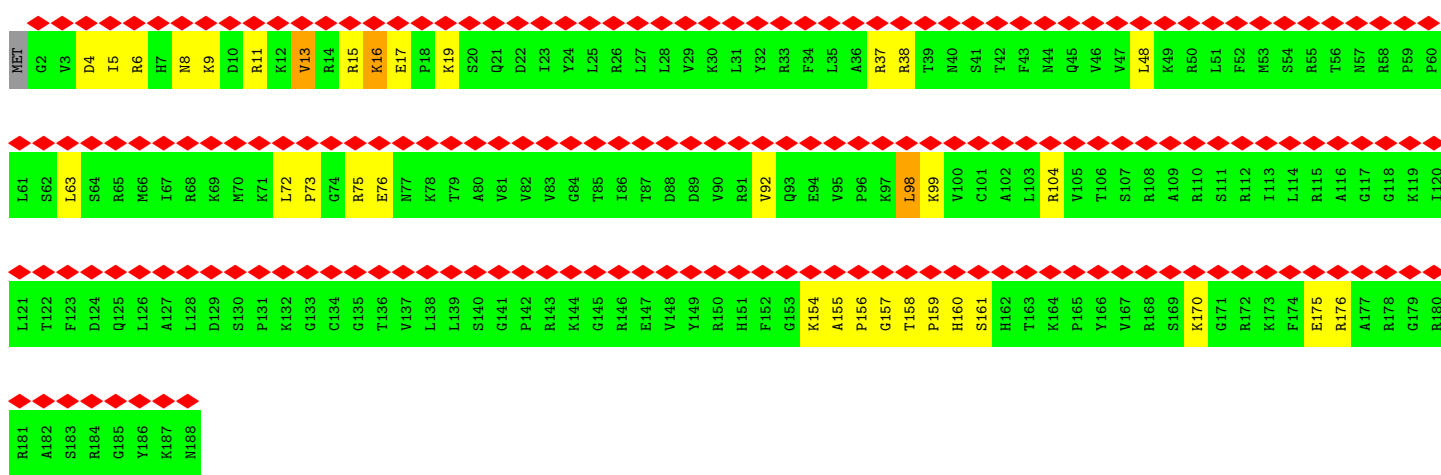
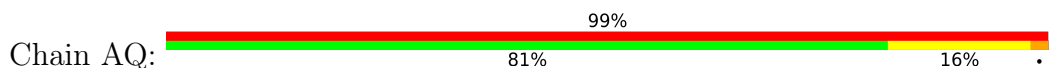
• Molecule 15: 60S ribosomal protein L14



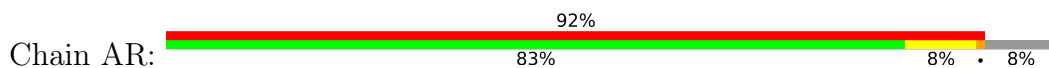
• Molecule 18: 60S ribosomal protein L17

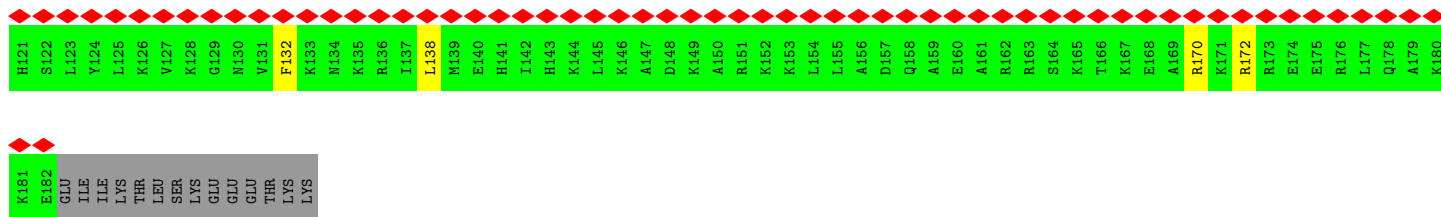


• Molecule 19: 60S ribosomal protein L18

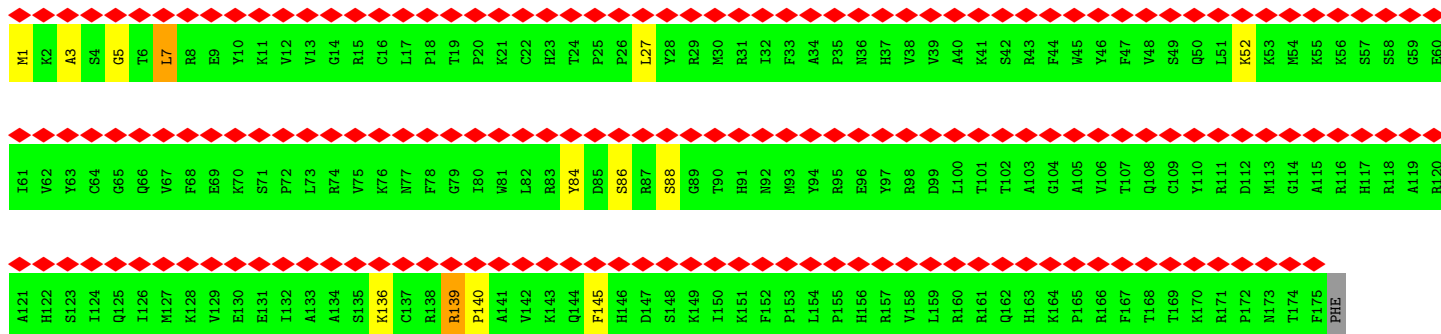
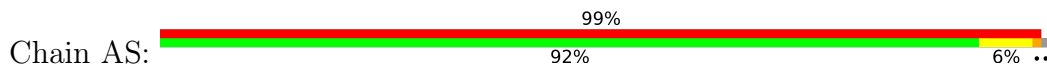


• Molecule 20: 60S ribosomal protein L19

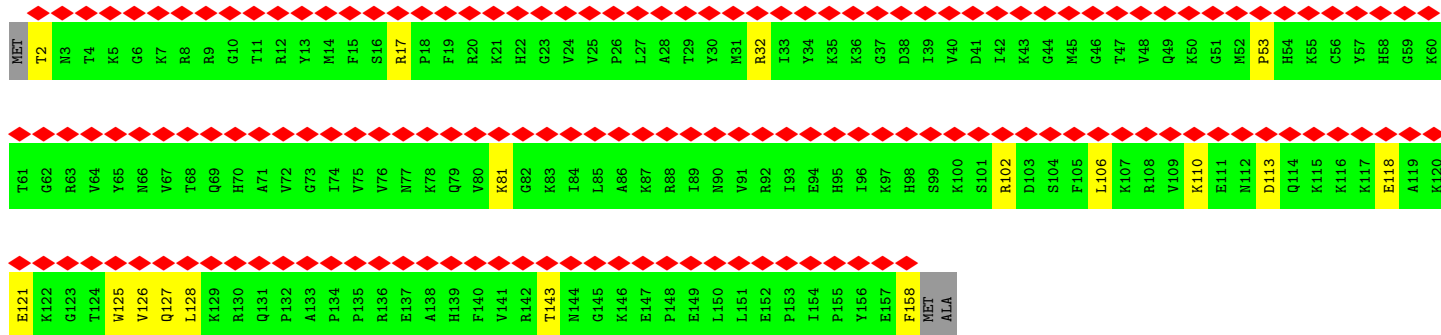
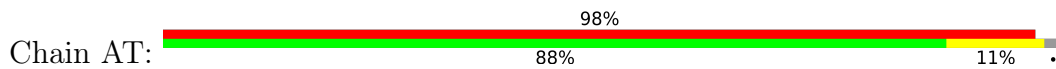




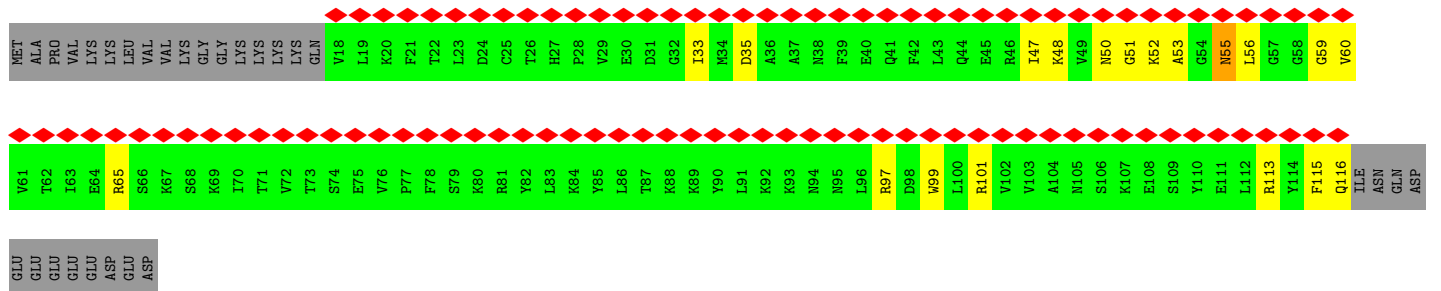
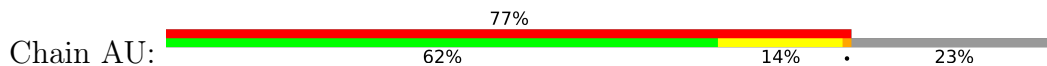
• Molecule 21: 60S ribosomal protein L18a

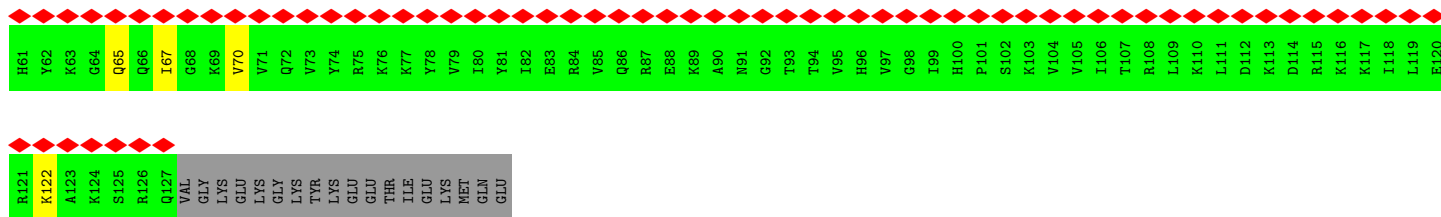


• Molecule 22: 60S ribosomal protein L21

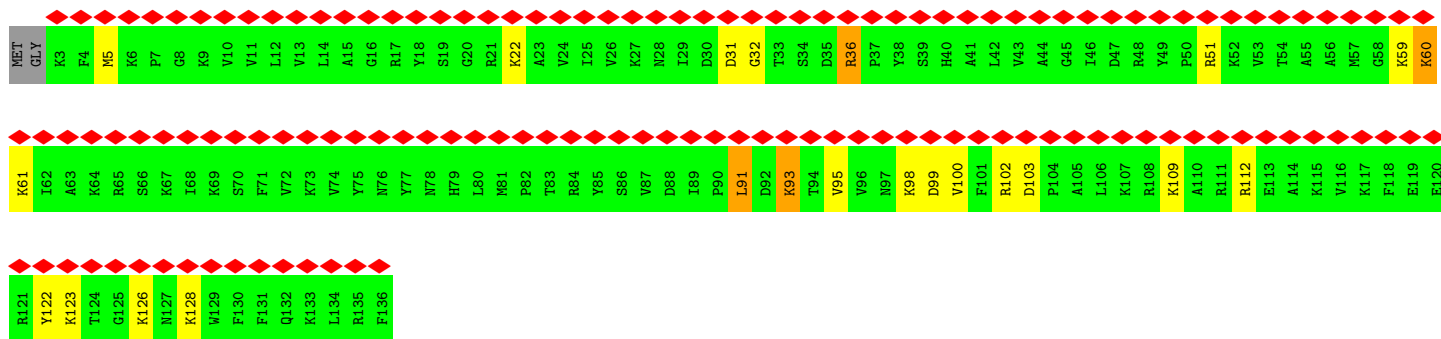
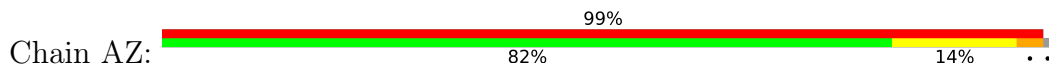


• Molecule 23: 60S ribosomal protein L22

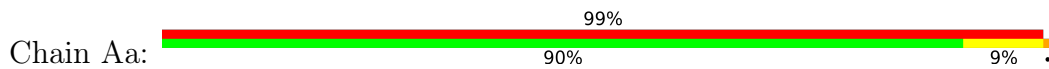




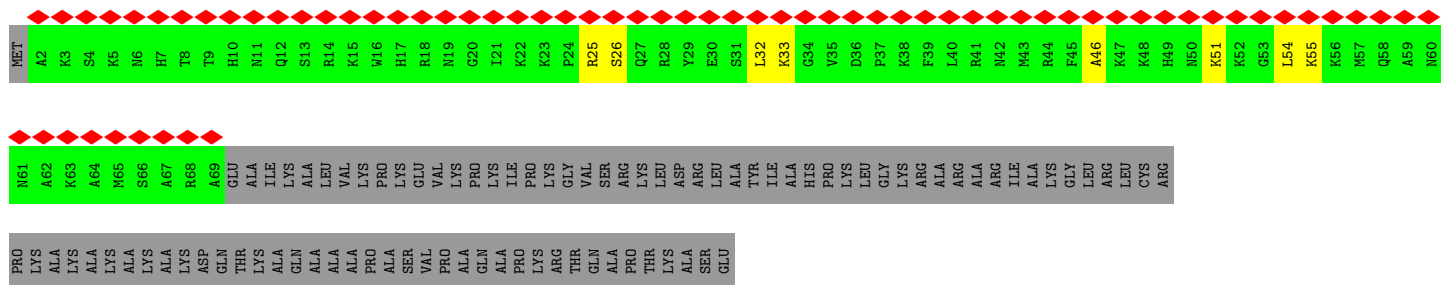
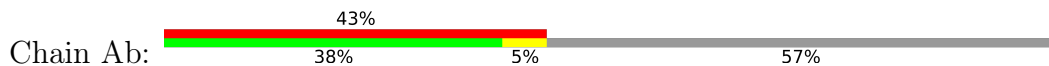
• Molecule 28: 60S ribosomal protein L27



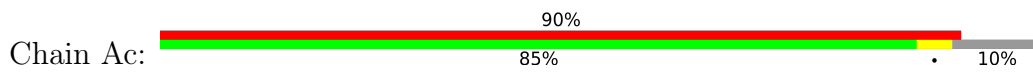
• Molecule 29: 60S ribosomal protein L27a



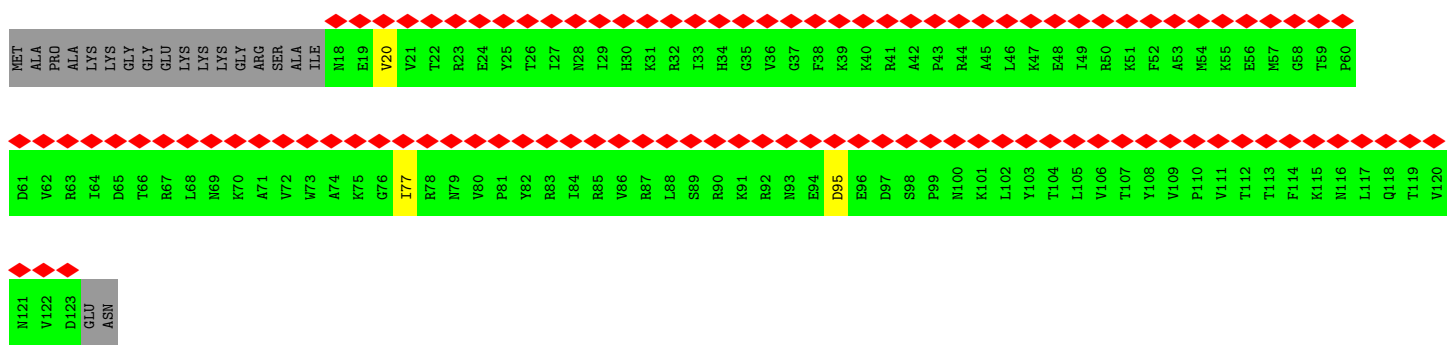
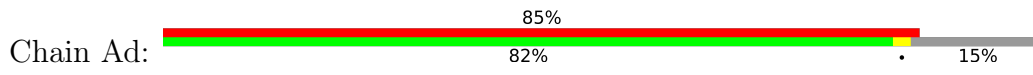
• Molecule 30: 60S ribosomal protein L29



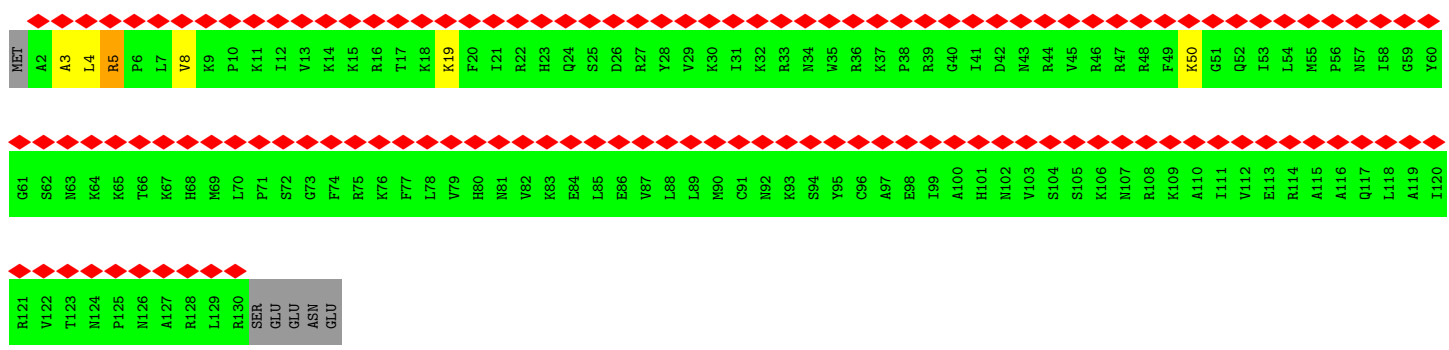
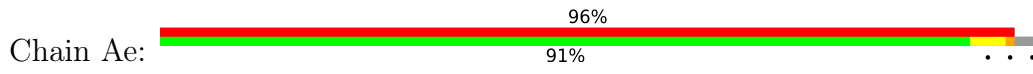
• Molecule 31: 60S ribosomal protein L30



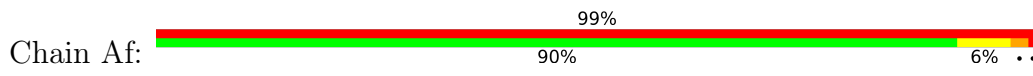
• Molecule 32: 60S ribosomal protein L31



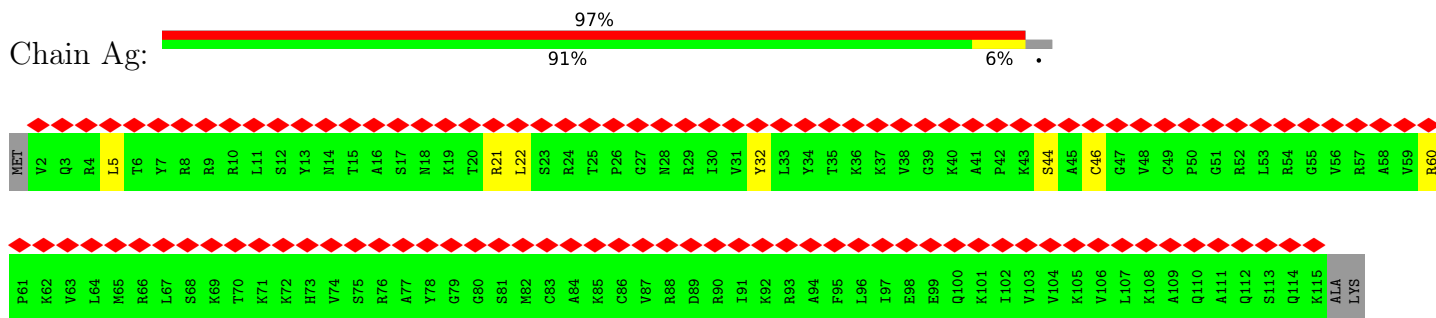
• Molecule 33: 60S ribosomal protein L32



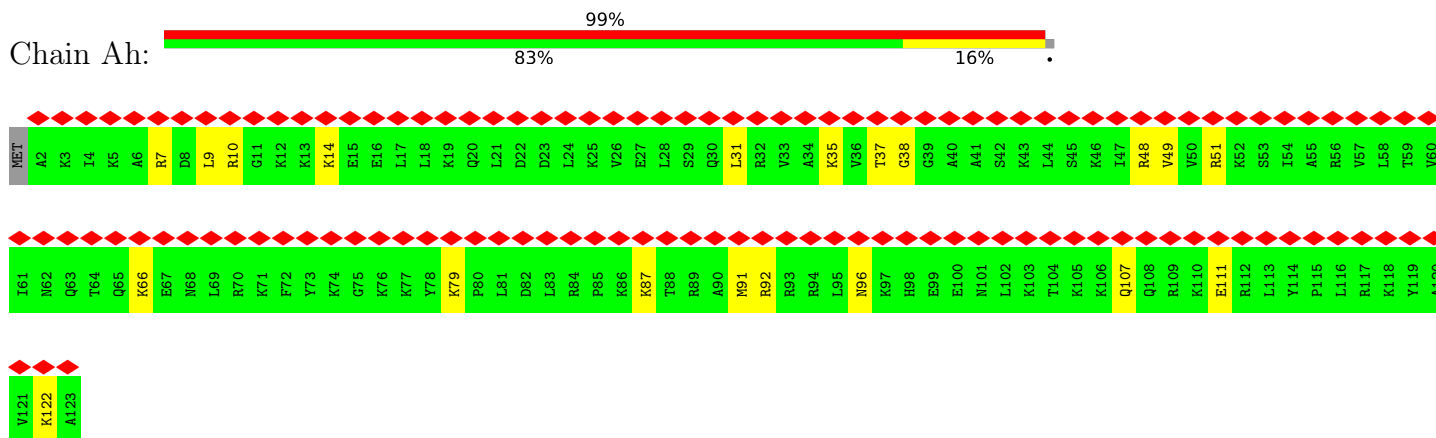
• Molecule 34: 60S ribosomal protein L35a



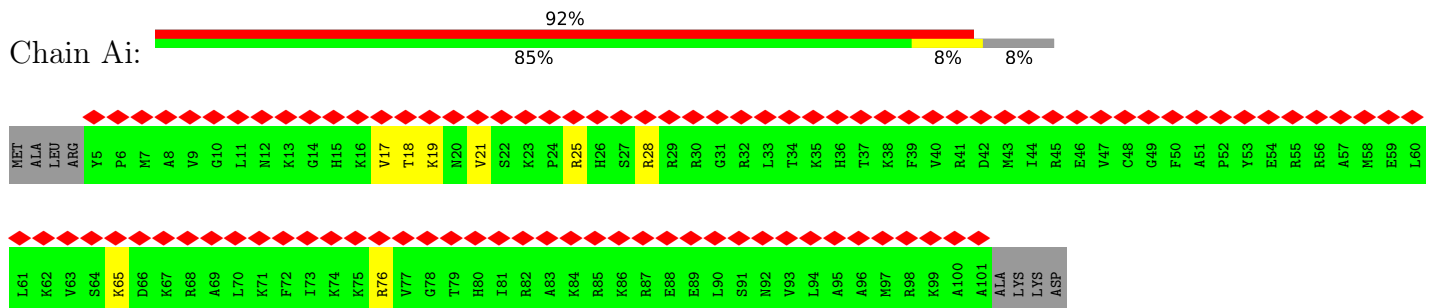
• Molecule 35: 60S ribosomal protein L34



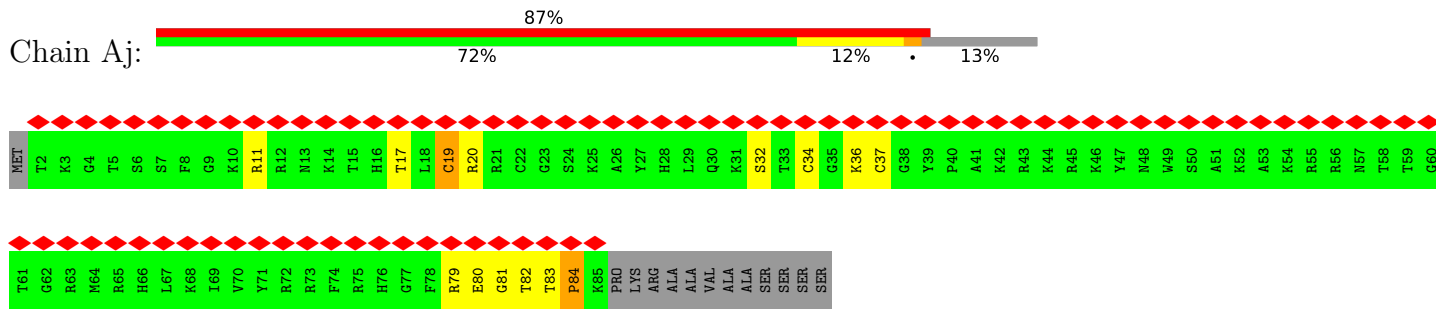
• Molecule 36: 60S ribosomal protein L35



• Molecule 37: 60S ribosomal protein L36



• Molecule 38: 60S ribosomal protein L37

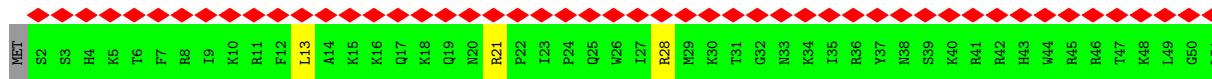
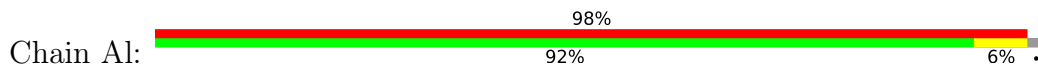


• Molecule 39: 60S ribosomal protein L38

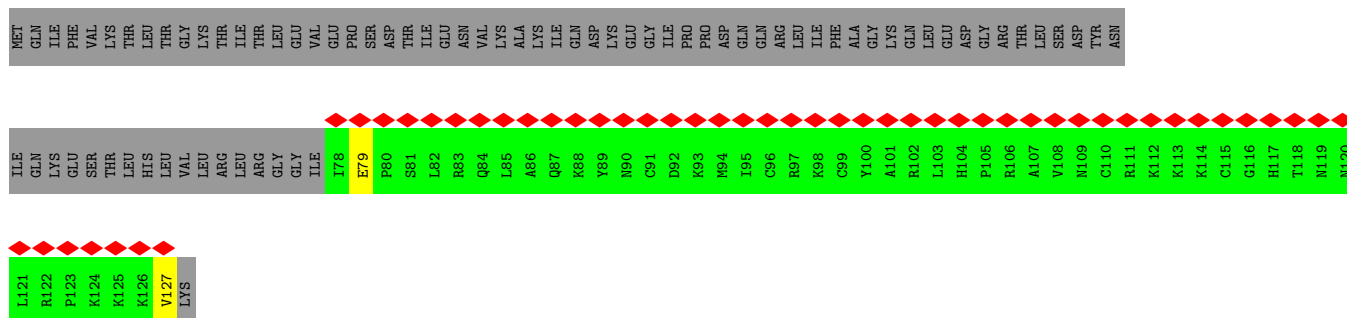




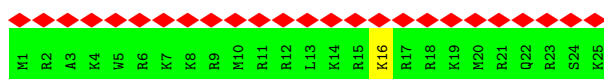
• Molecule 40: 60S ribosomal protein L39



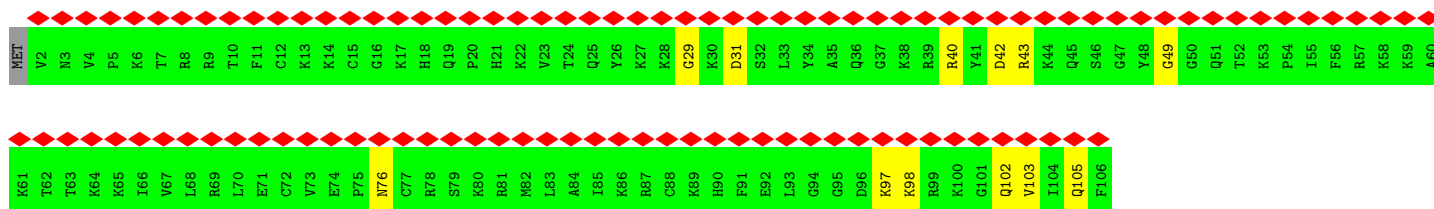
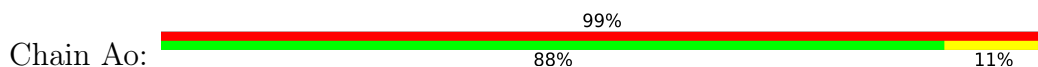
• Molecule 41: Ubiquitin-60S ribosomal protein L40



• Molecule 42: 60S ribosomal protein L41

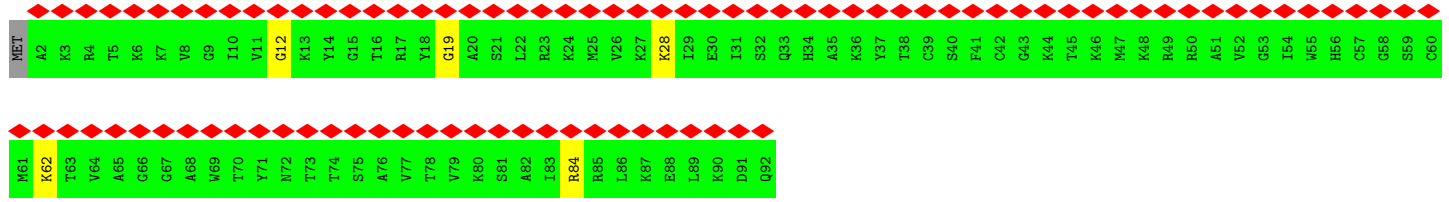


• Molecule 43: 60S ribosomal protein L36a

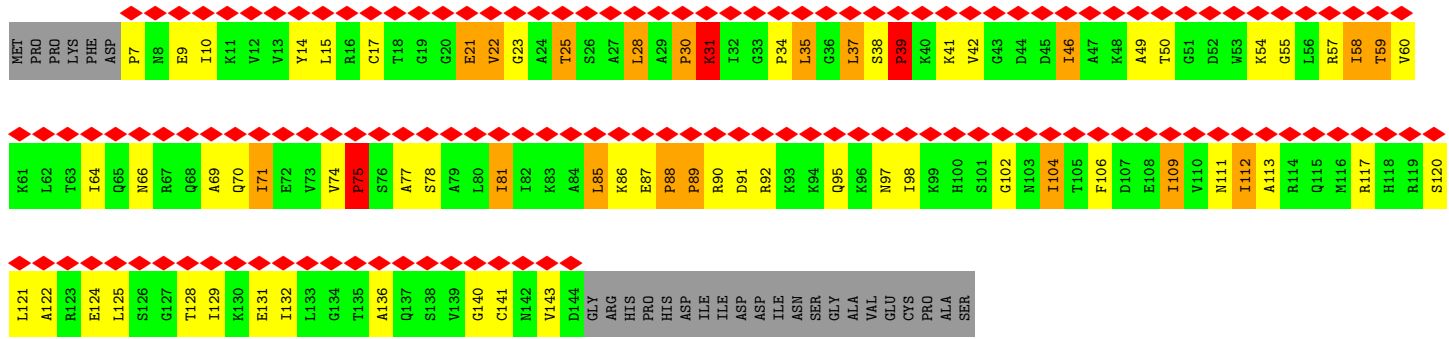
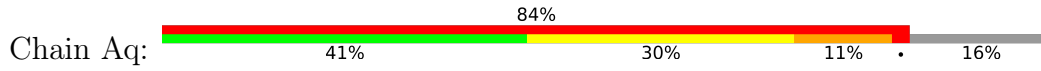


• Molecule 44: 60S ribosomal protein L37a

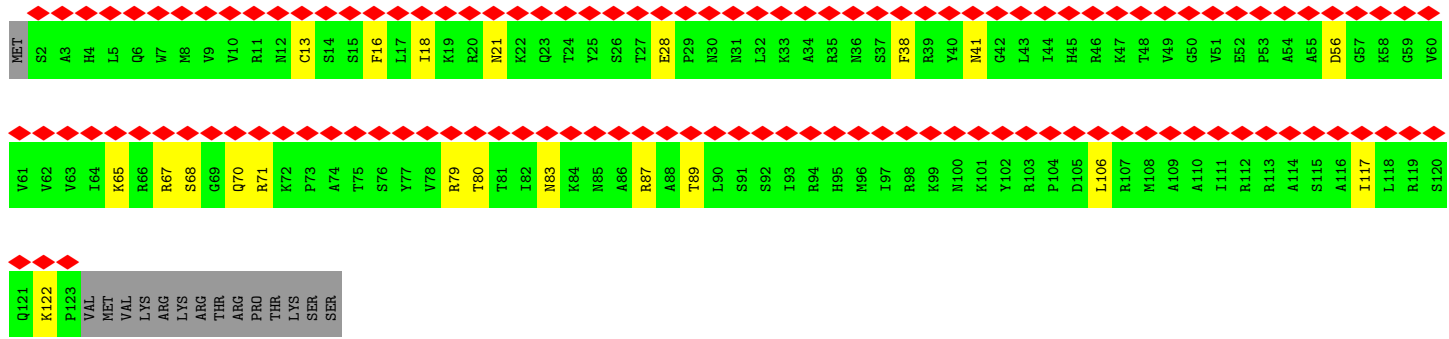
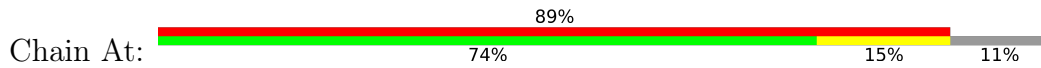




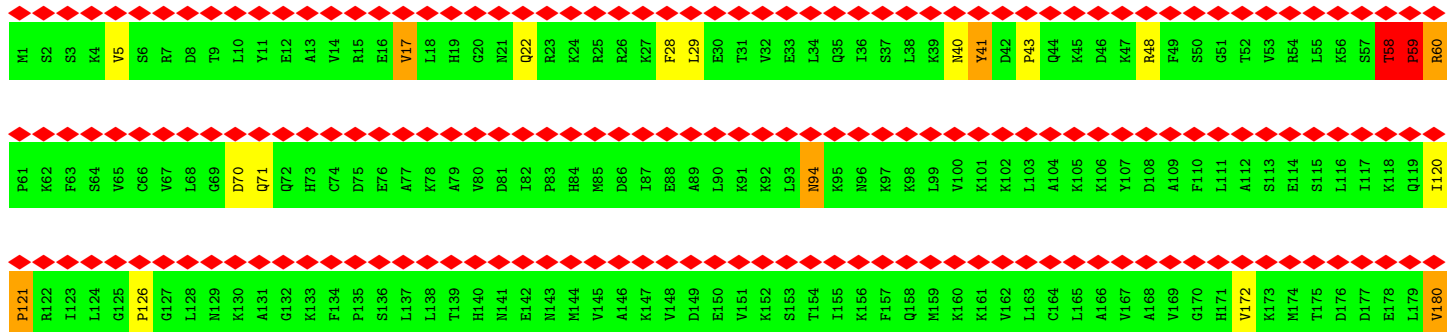
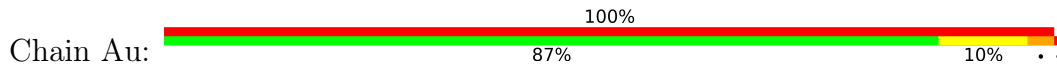
• Molecule 45: 60S ribosomal protein L12

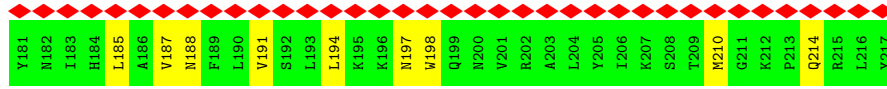


• Molecule 46: 60S ribosomal protein L28

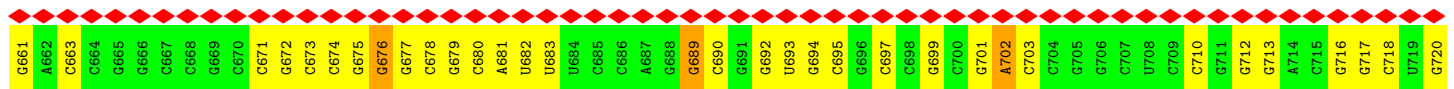
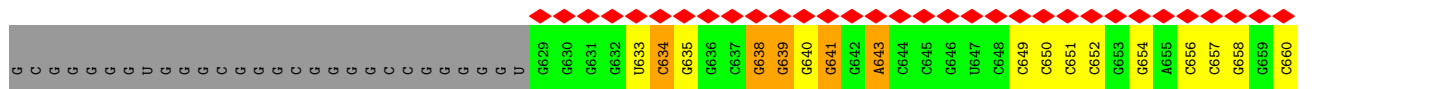
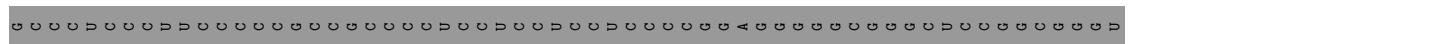
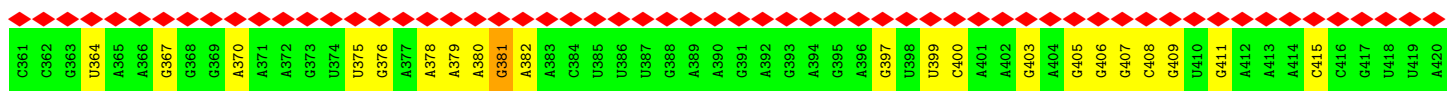
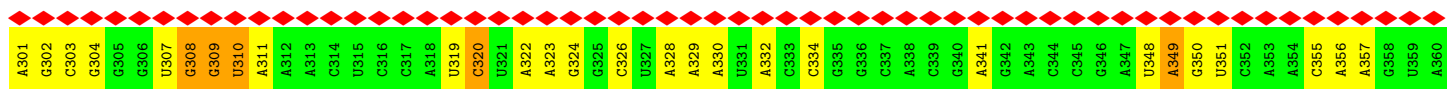
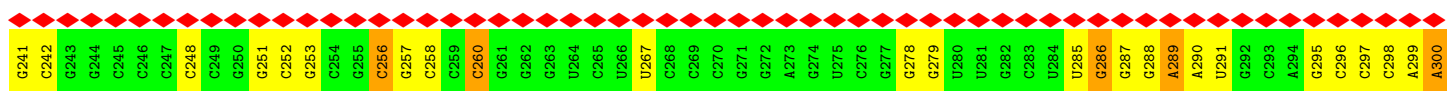
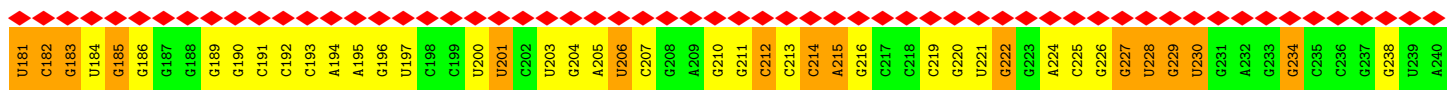
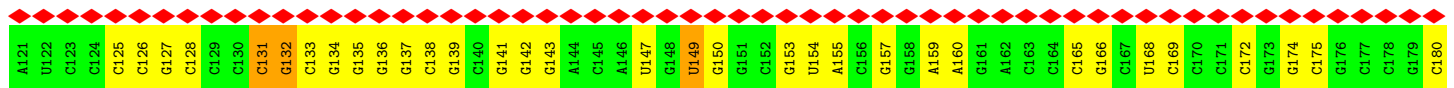
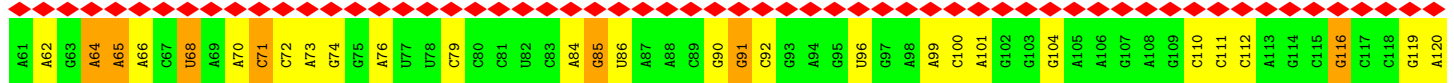
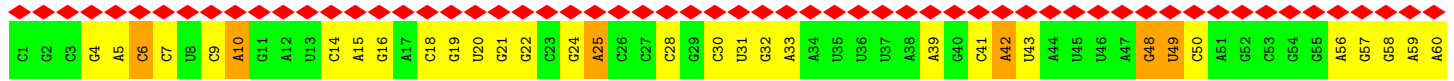
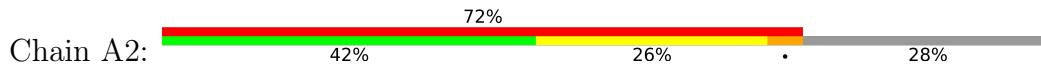


• Molecule 47: 60S ribosomal protein L10a



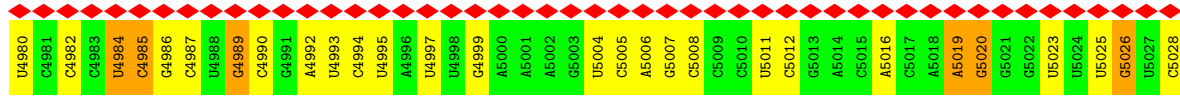


• Molecule 48: 28S ribosomal RNA

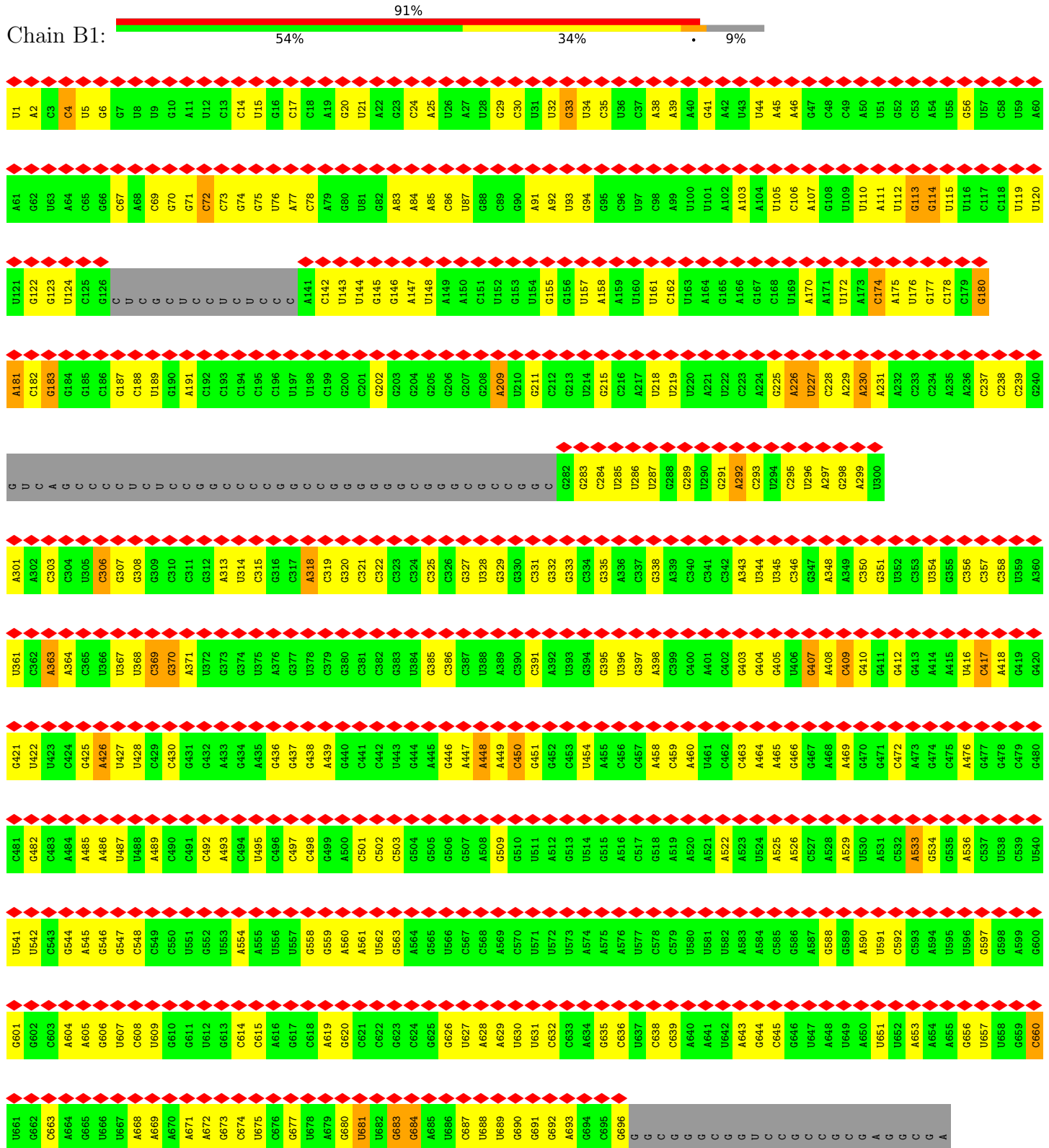


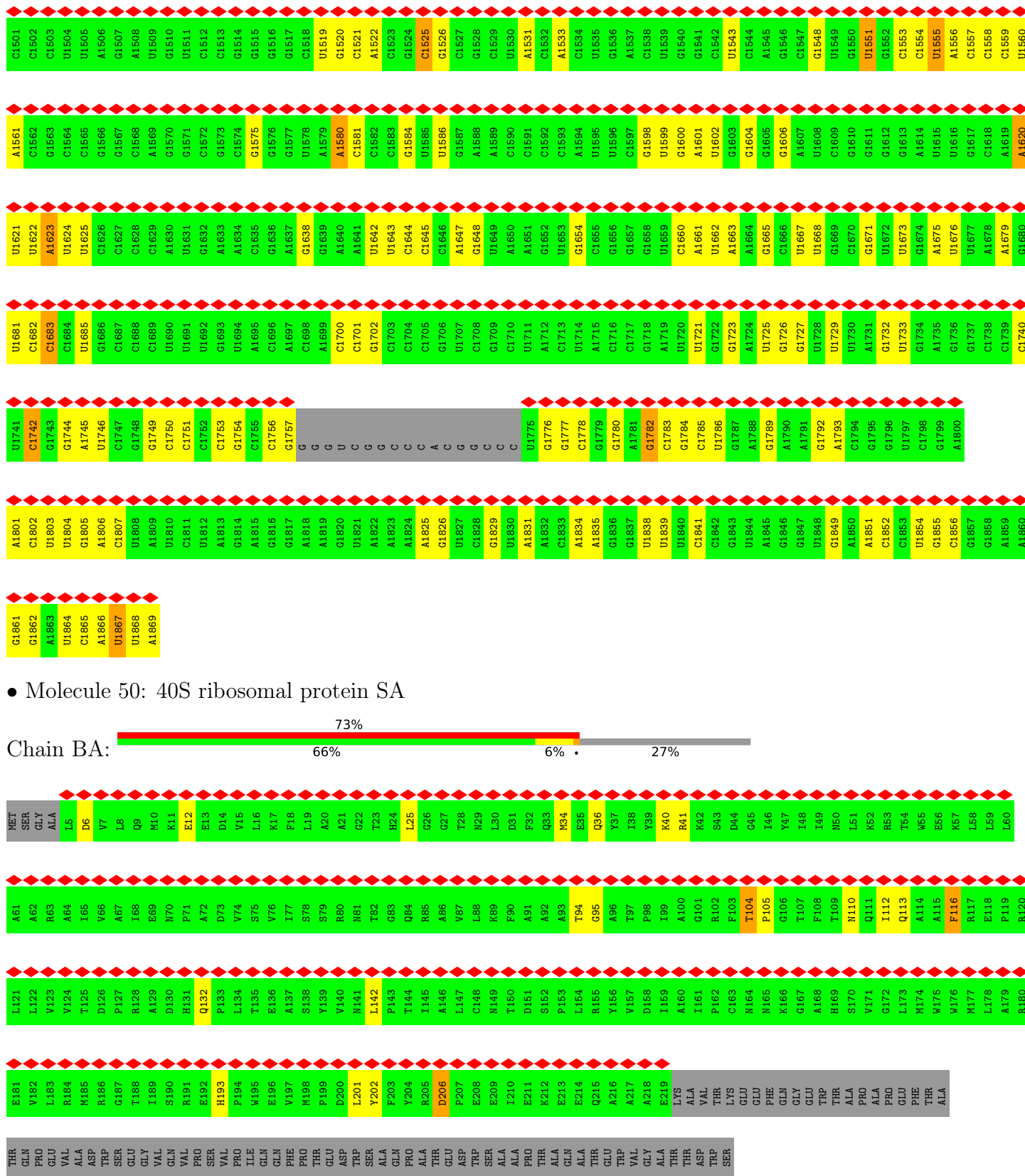
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U2283	C2043	G1803	G1743	G1623	C1503
U2284	G2044	U1804	C1744	A1624	G1504
G2285	G2045	G1805	C1745	C1625	A1505
A2286	G2046	G1806	G1746	C1626	A1506
A2287	G2047	G1807	A1747	C1627	A1507
G2288	C2048	G1808	A1748	A1628	G1508
G2289	G2049	C1809	A1749	A1629	A1509
C2290	A2050	C1810	C1750	C1630	U1510
U2291	U2051	G1811	G1751	U1631	G1511
A2292	A2052	G1812	A1752	A1632	G1512
G2293	C2053	G1813	U1753	G1633	U1513
G2294	A1993	G1814	C1754	G1633	G1514
G2295	A1994	G1815	G1755	U1635	A1515
G2296	C1995	U1815	U1756	G1636	A1516
G2297	U1996	G1816	C1756	G1636	C1517
G2298	G1997	G1817	A1757	C1637	U1518
G2299	A1998	G1818	A1758	U1638	A1519
G2300	C1999	A1819	C1759	U1639	U1520
G2301	C2000	U1820	G1760	G1640	G1521
C2302	U2001	G1821	U1761	U1641	G1522
C2303	G2002	G1822	A1762	U1642	C1523
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G2305	C2004	A1824	U1764	C1644	G1525
G2306	G2005	U1825	U1765	U1645	G1526
G2307	A2006	G1826	U1766	U1645	G1527
U2308	A2007	U1827	C1767	C1647	C1528
G2309	U2008	G1828	A1768	C1648	A1529
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G2312	C1952	G1832	U1772	G1652	C1533
A2313	G1953	U1833	U1773	U1653	G1534
G2314	G1954	G1834	U1774	U1654	A1535
G2315	G2015	G1835	A1775	U1655	A1536
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G2318	C2018	G1838	U1778	C1658	C1539
G2319	U2019	A1839	G1779	U1659	A1540
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G2322	A2022	U1842	U1782	G1662	G1543
C2323	A2023	U1843	A1783	G1663	G1544
G2324	A2024	A1904	U1784	A1664	A1545
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A2326	G1966	G1846	A1786	U1666	A1547
G2327	G2026	U1847	A1787	G1667	C1548
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G2330	U2029	G1909	U1790	G1670	U1551
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U2332	G2031	U1911	C1791	G1672	A1613
G2333	C2032	A1912	U1792	C1673	U1553
G2334	G2033	G1913	G1793	C1674	A1614
U2335	C2034	A1914	U1794	U1675	G1615
G2336	U2035	A1915	U1795	C1676	A1616
G2337	G2036	C1916	C1796	U1677	C1617
U2338	G2037	C1917	G1797	C1678	A1558
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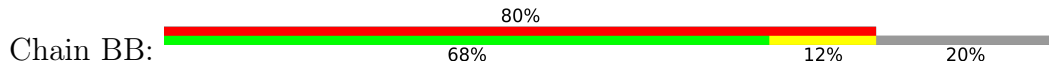
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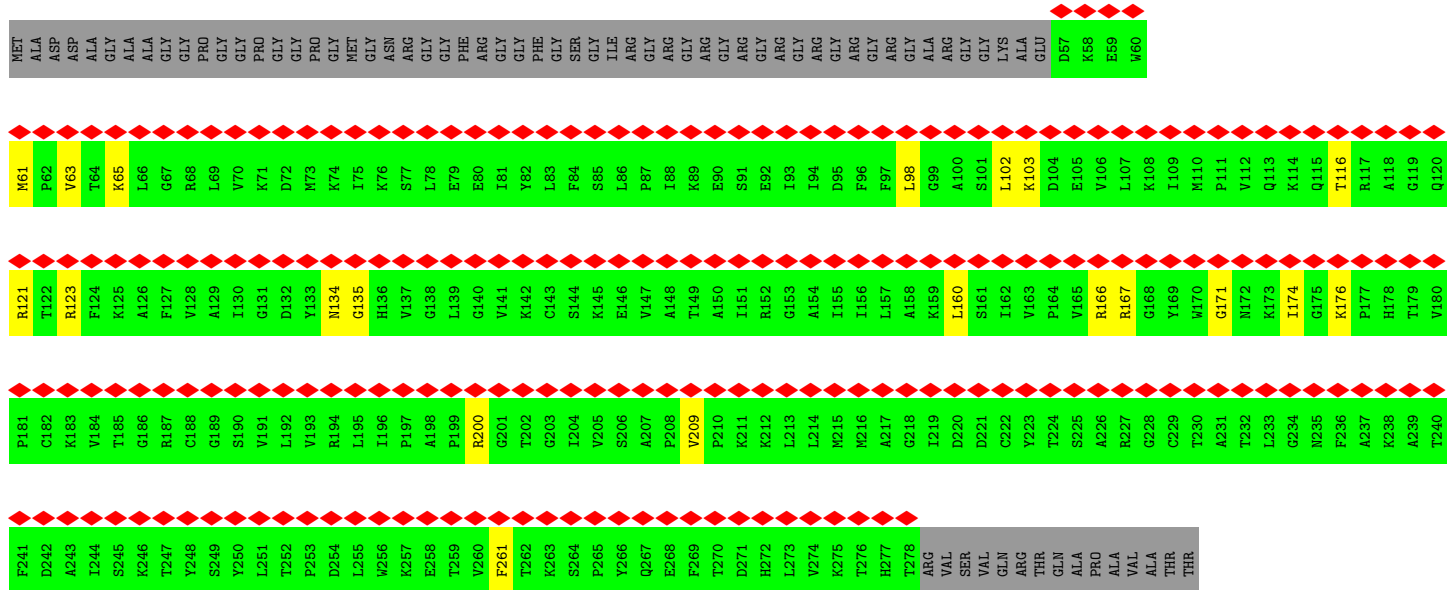
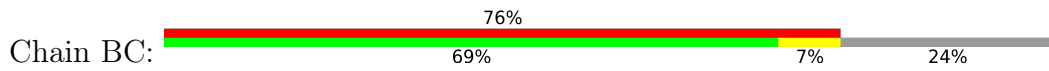
● Molecule 49: 18S ribosomal RNA



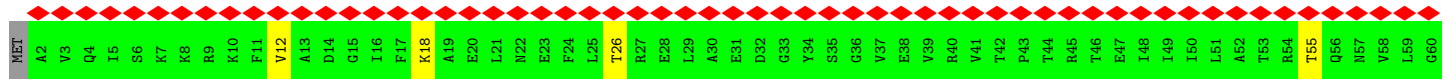
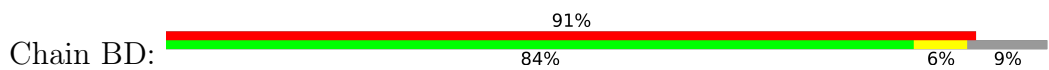


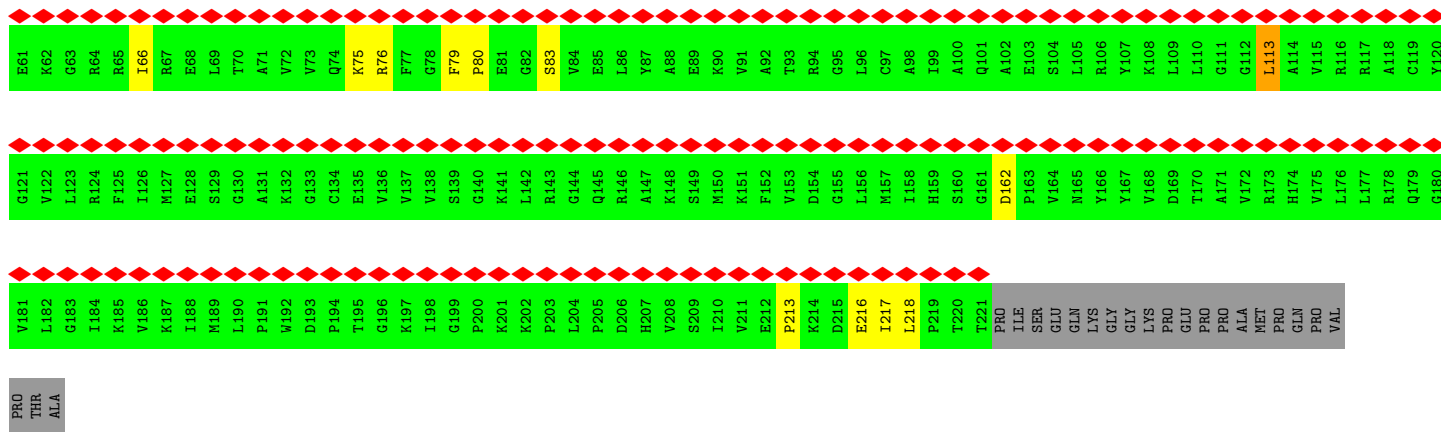


• Molecule 52: 40S ribosomal protein S2

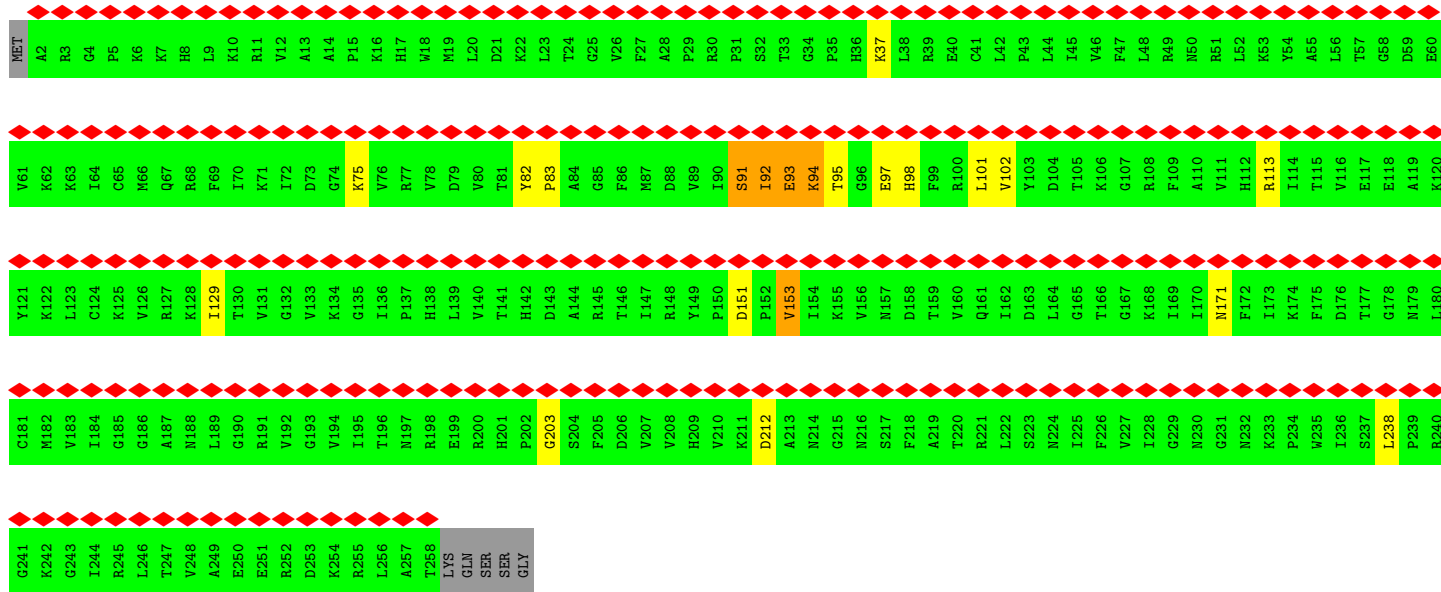
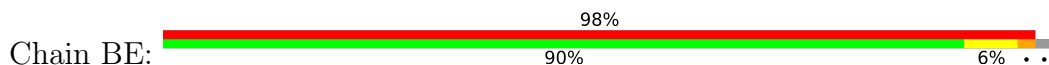


• Molecule 53: 40S ribosomal protein S3

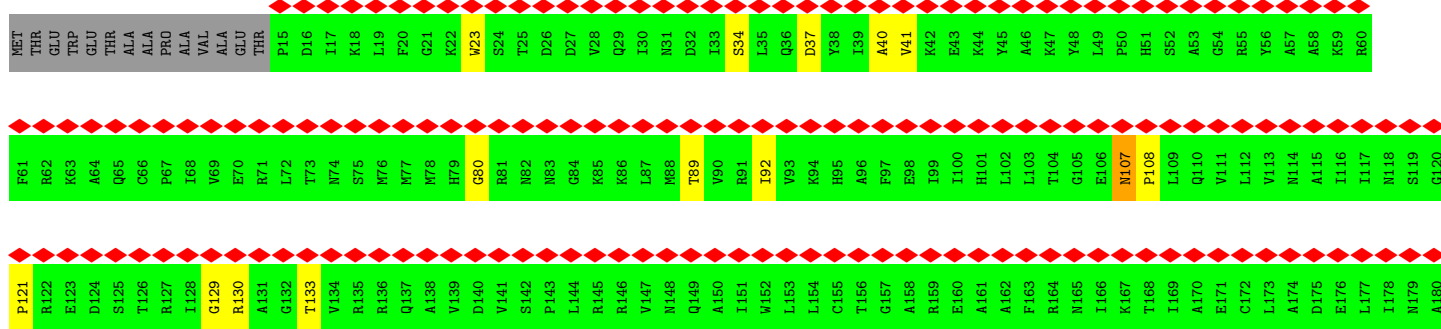
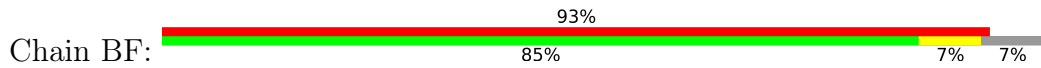


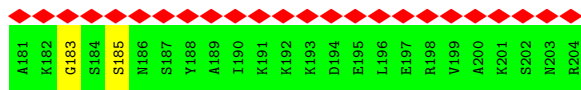


• Molecule 54: 40S ribosomal protein S4, Y isoform 1

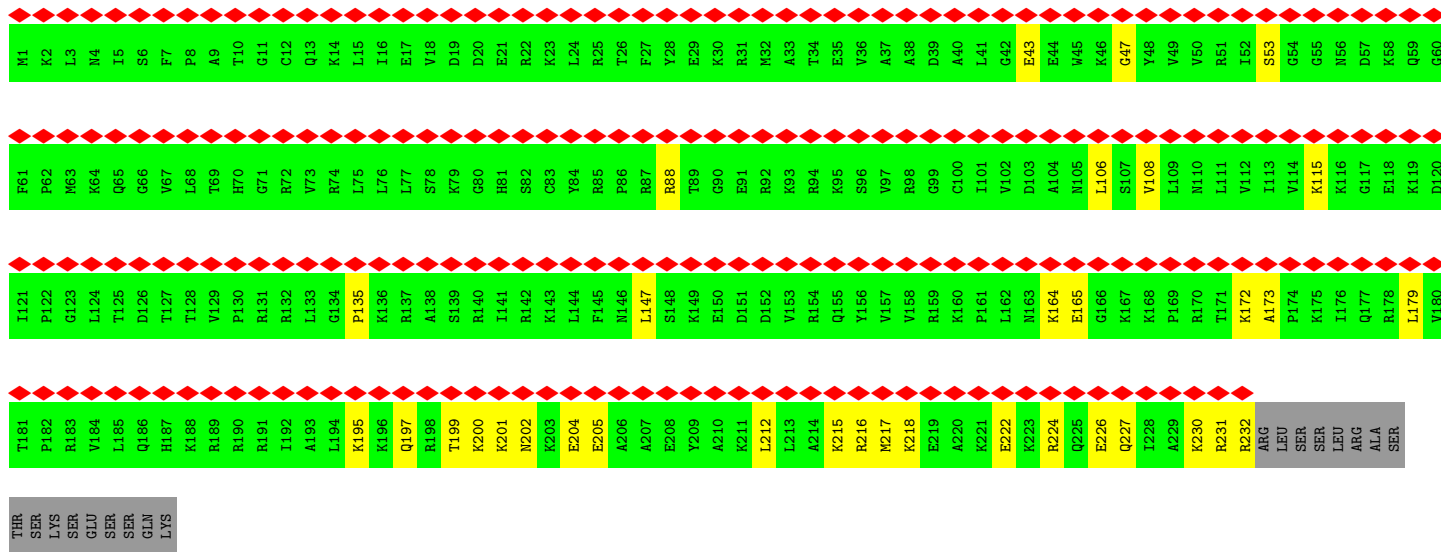
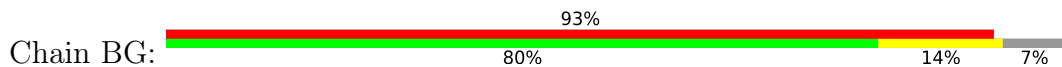


• Molecule 55: 40S ribosomal protein S5

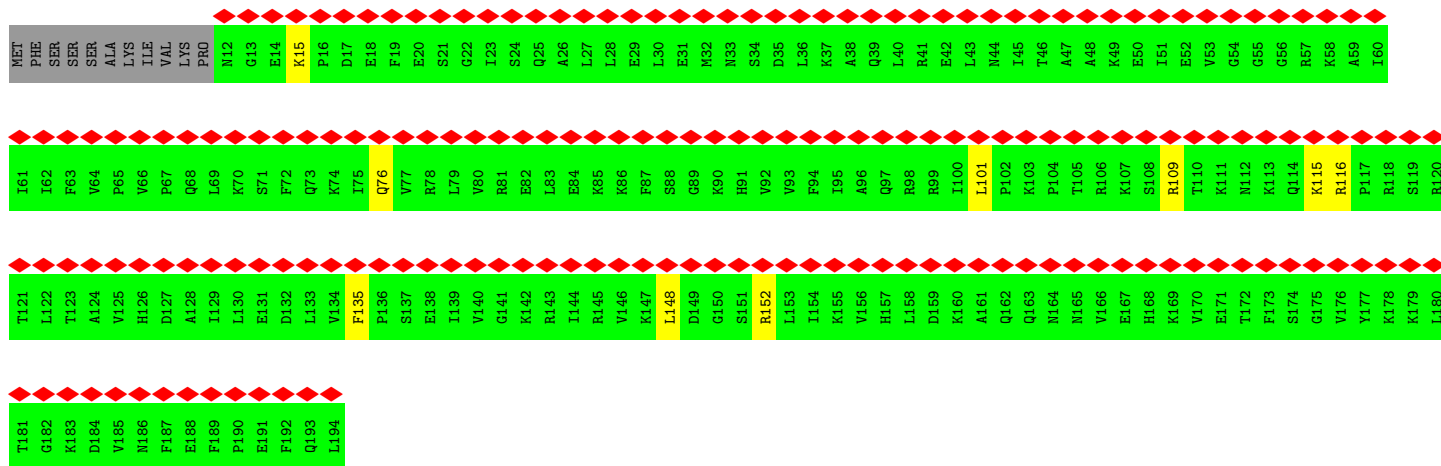
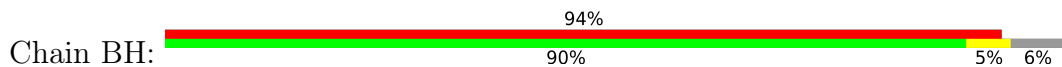




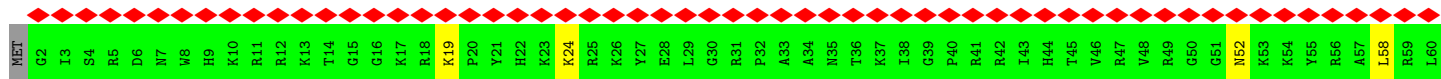
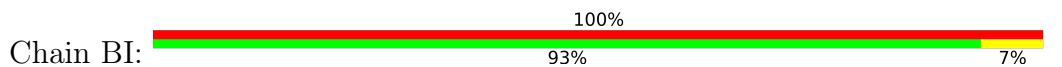
• Molecule 56: 40S ribosomal protein S6



• Molecule 57: 40S ribosomal protein S7

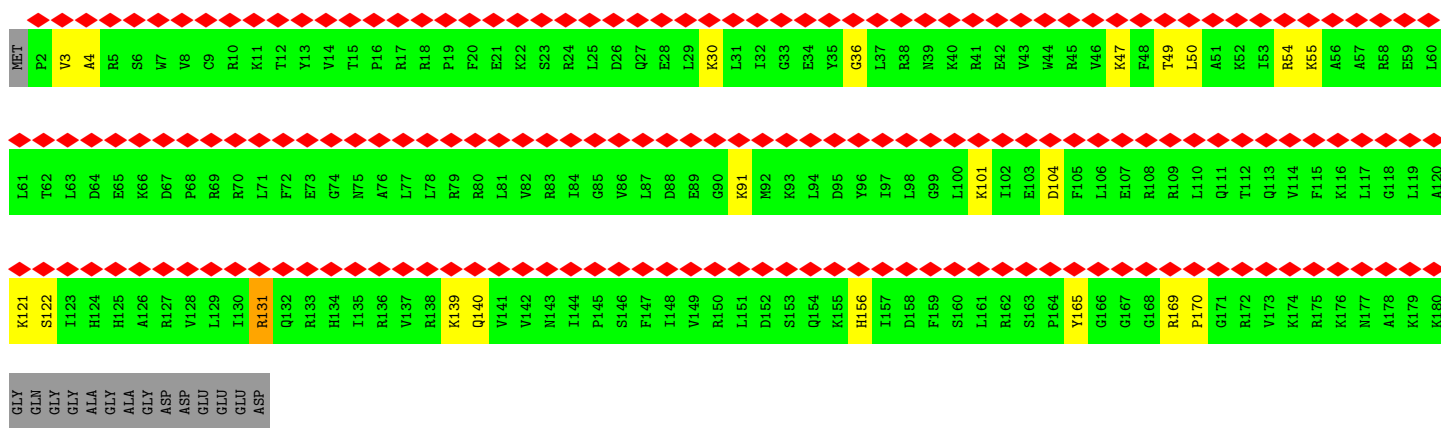
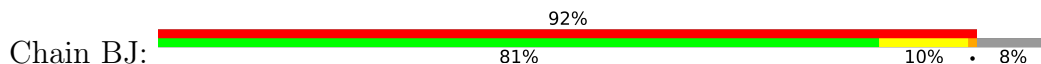


• Molecule 58: 40S ribosomal protein S8

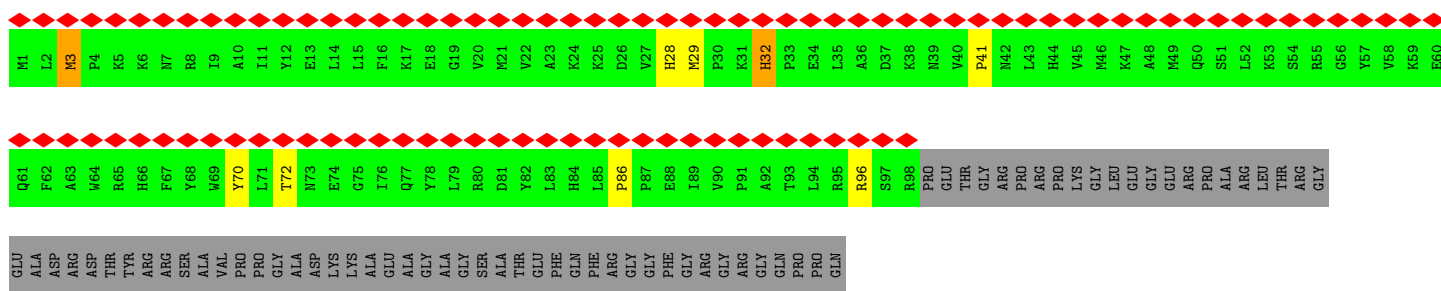




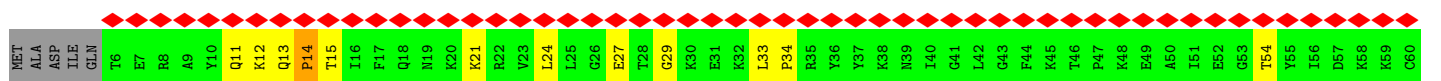
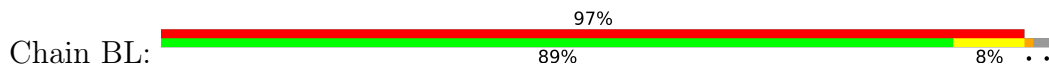
• Molecule 59: 40S ribosomal protein S9

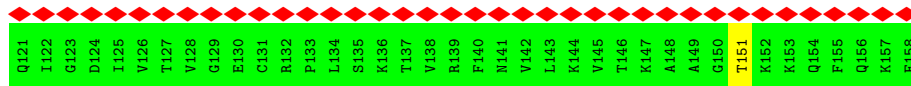
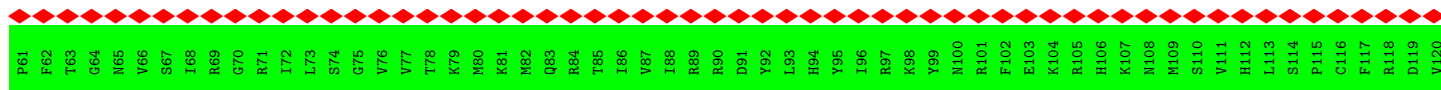


• Molecule 60: 40S ribosomal protein S10

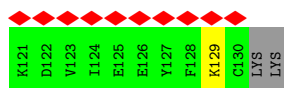
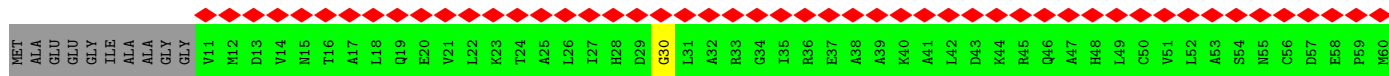
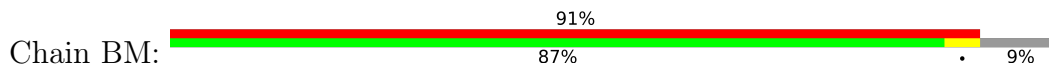


• Molecule 61: 40S ribosomal protein S11

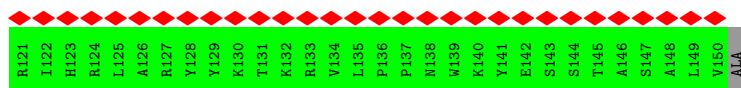
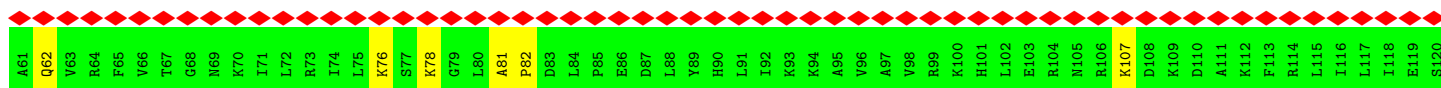
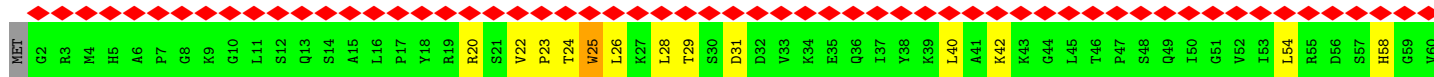
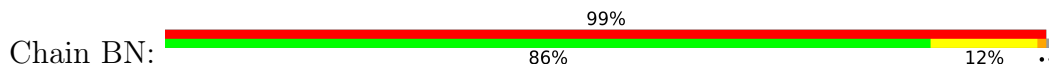




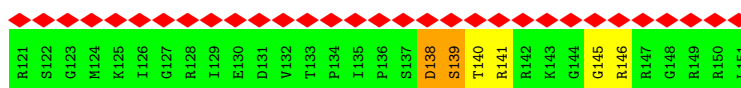
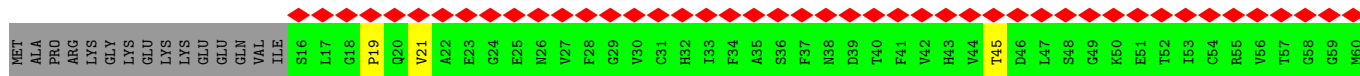
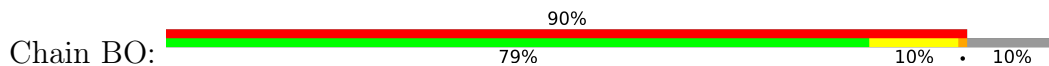
• Molecule 62: 40S ribosomal protein S12



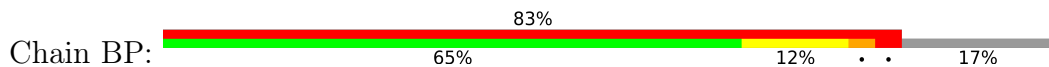
• Molecule 63: 40S ribosomal protein S13



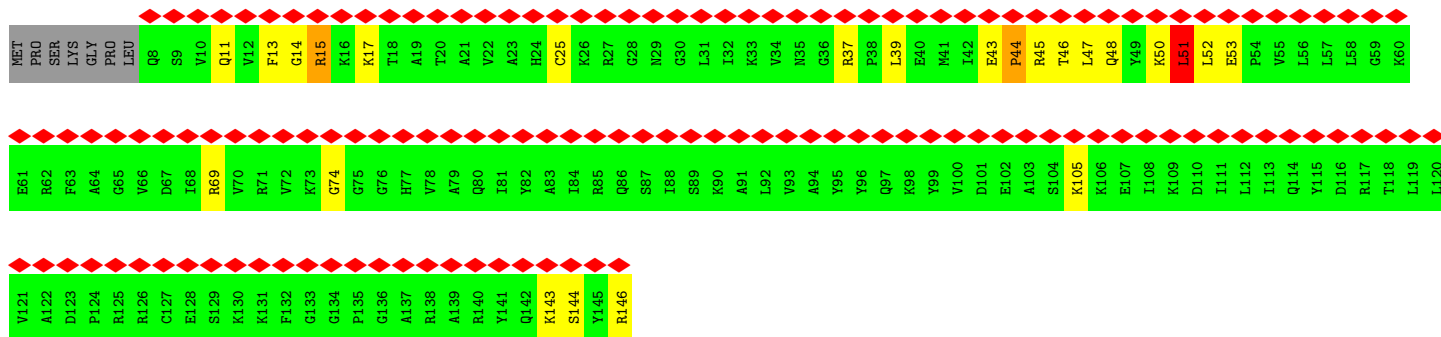
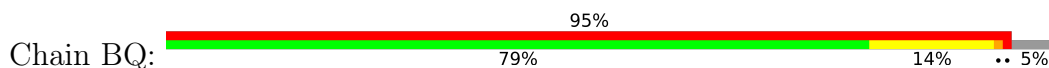
• Molecule 64: 40S ribosomal protein S14



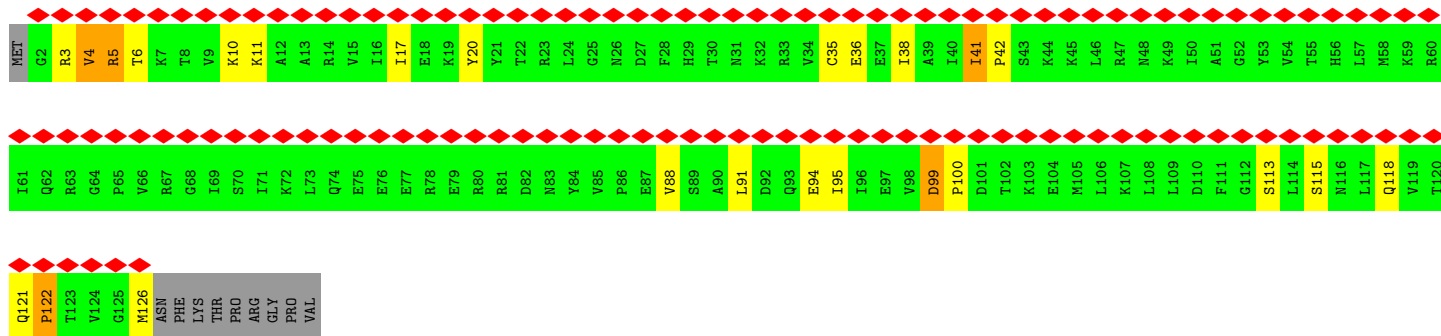
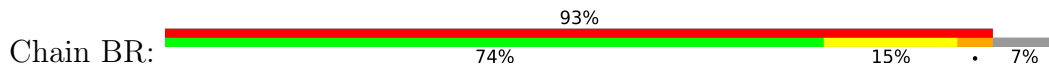
• Molecule 65: 40S ribosomal protein S15



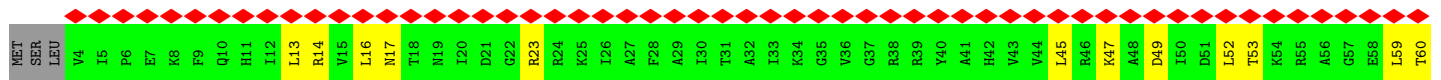
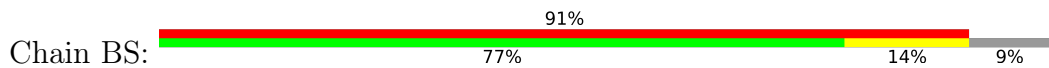
• Molecule 66: 40S ribosomal protein S16

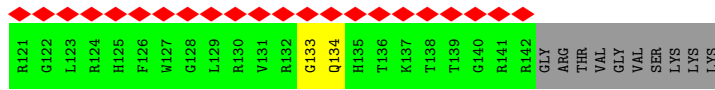
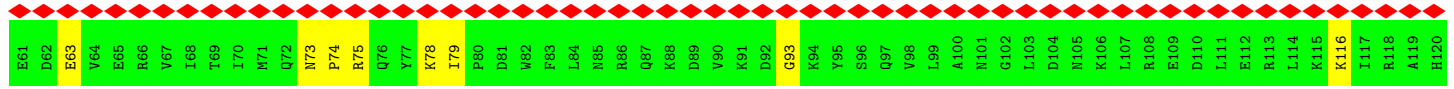


• Molecule 67: 40S ribosomal protein S17-like

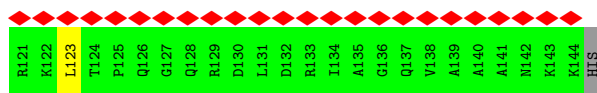
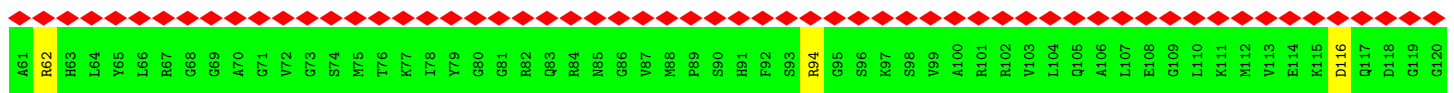
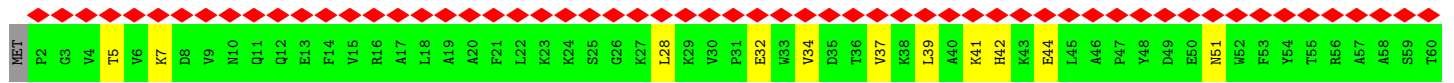
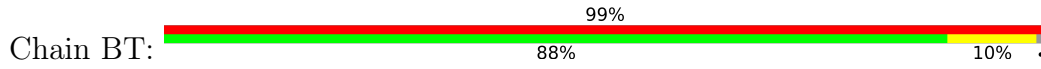


• Molecule 68: 40S ribosomal protein S18

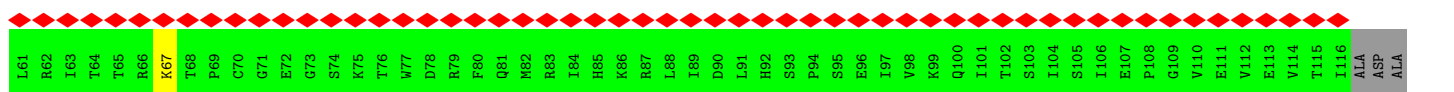
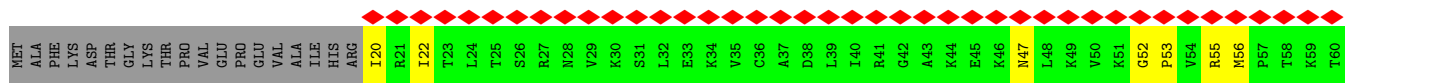
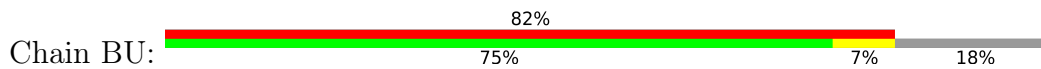




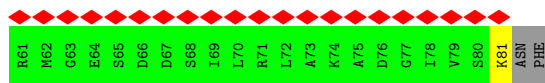
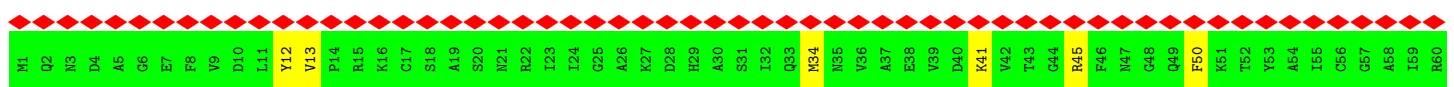
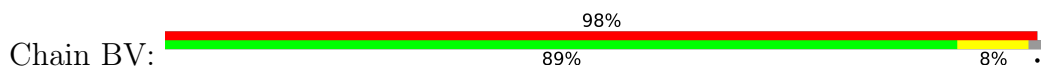
• Molecule 69: 40S ribosomal protein S19



• Molecule 70: 40S ribosomal protein S20

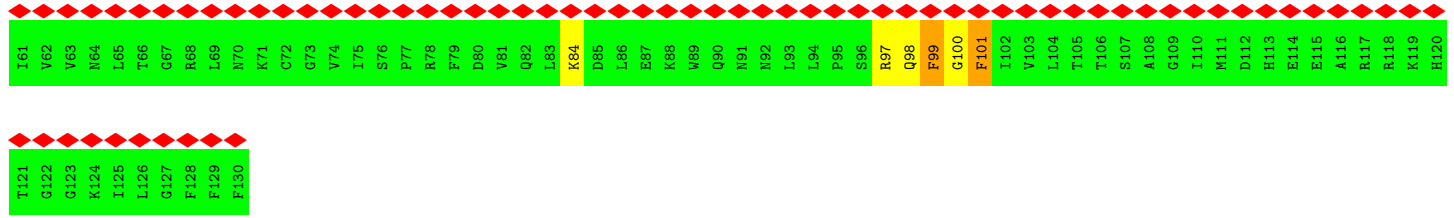


• Molecule 71: 40S ribosomal protein S21

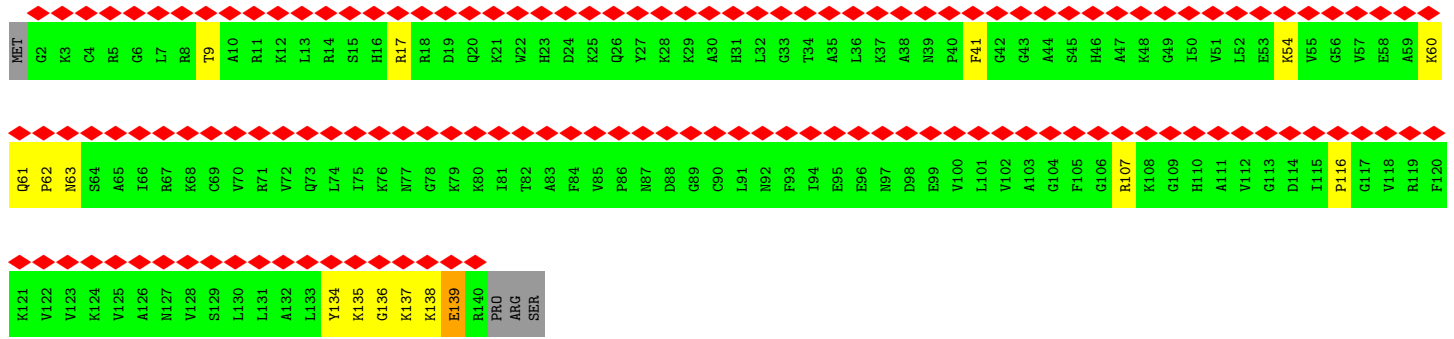
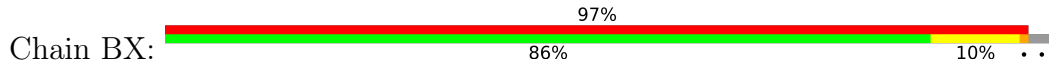


• Molecule 72: 40S ribosomal protein S15a

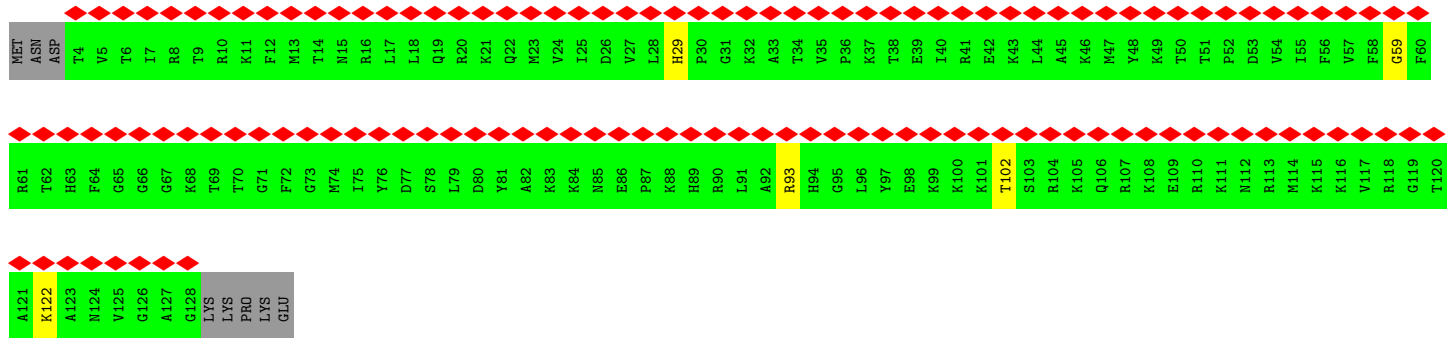
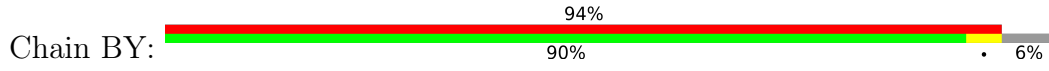




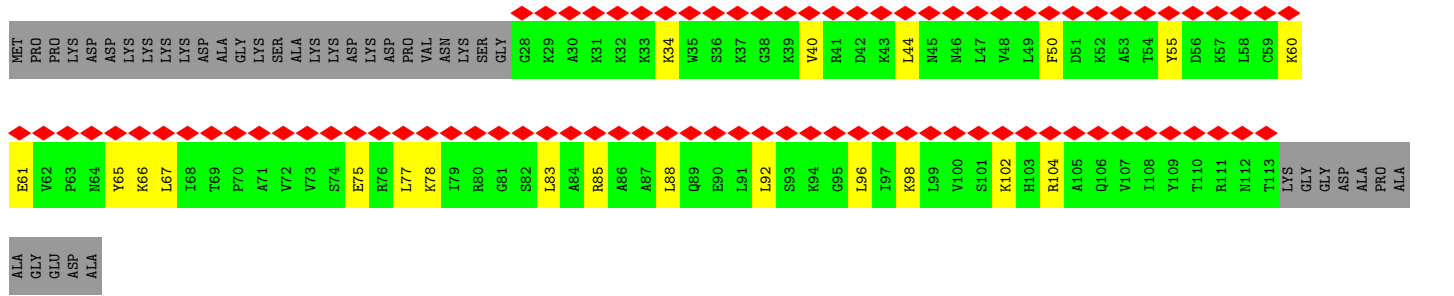
• Molecule 73: 40S ribosomal protein S23



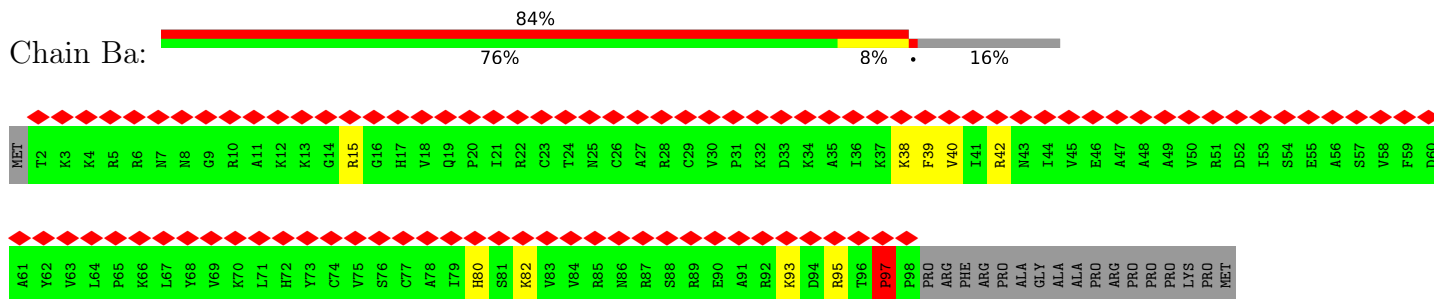
• Molecule 74: 40S ribosomal protein S24



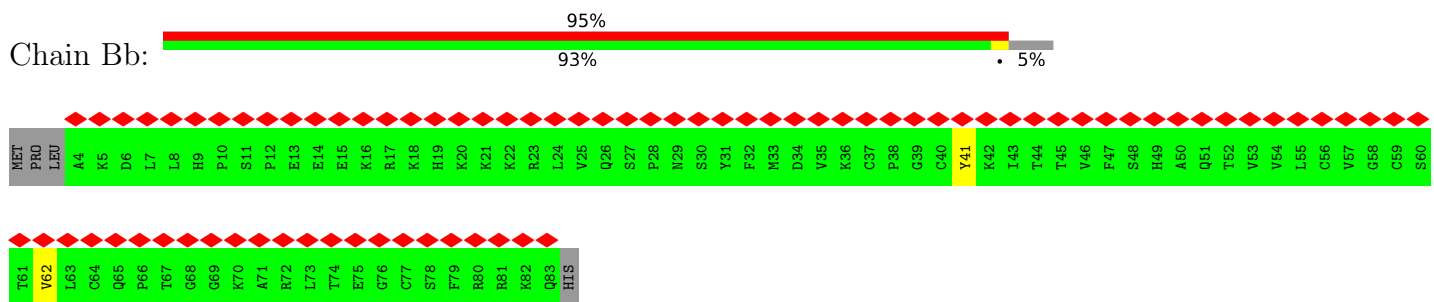
• Molecule 75: 40S ribosomal protein S25



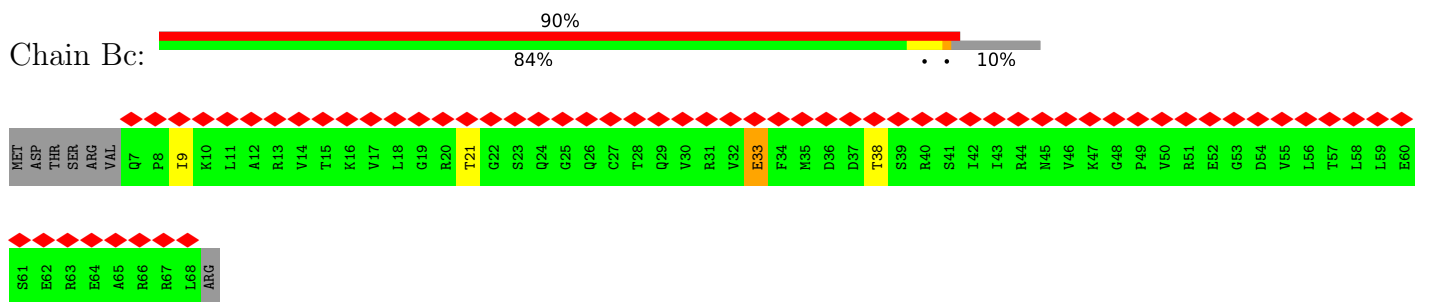
- Molecule 76: 40S ribosomal protein S26



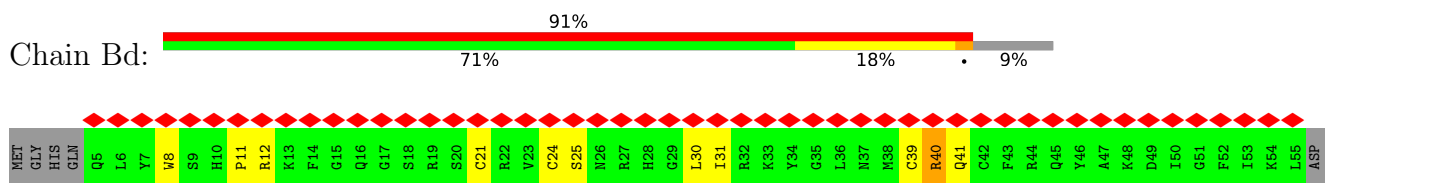
- Molecule 77: 40S ribosomal protein S27



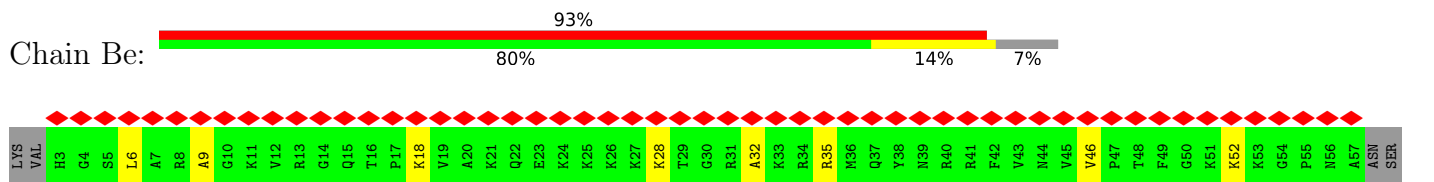
- Molecule 78: 40S ribosomal protein S28



- Molecule 79: 40S ribosomal protein S29

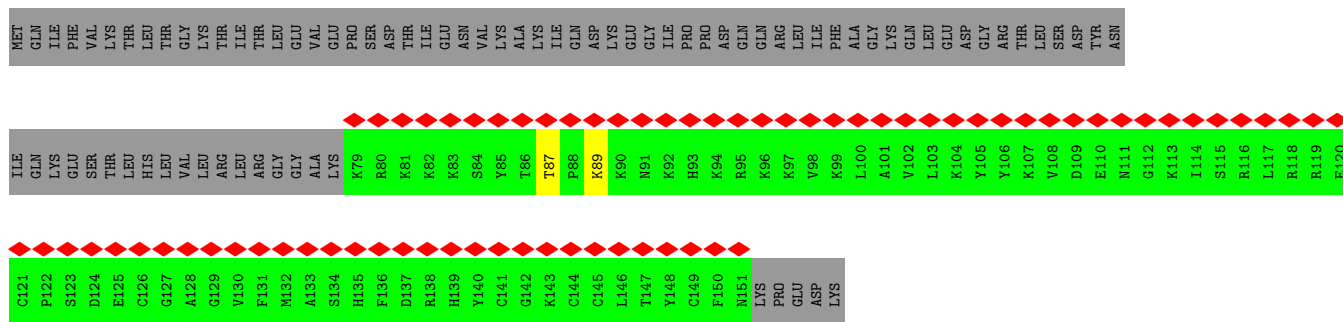


- Molecule 80: 40S ribosomal protein S30

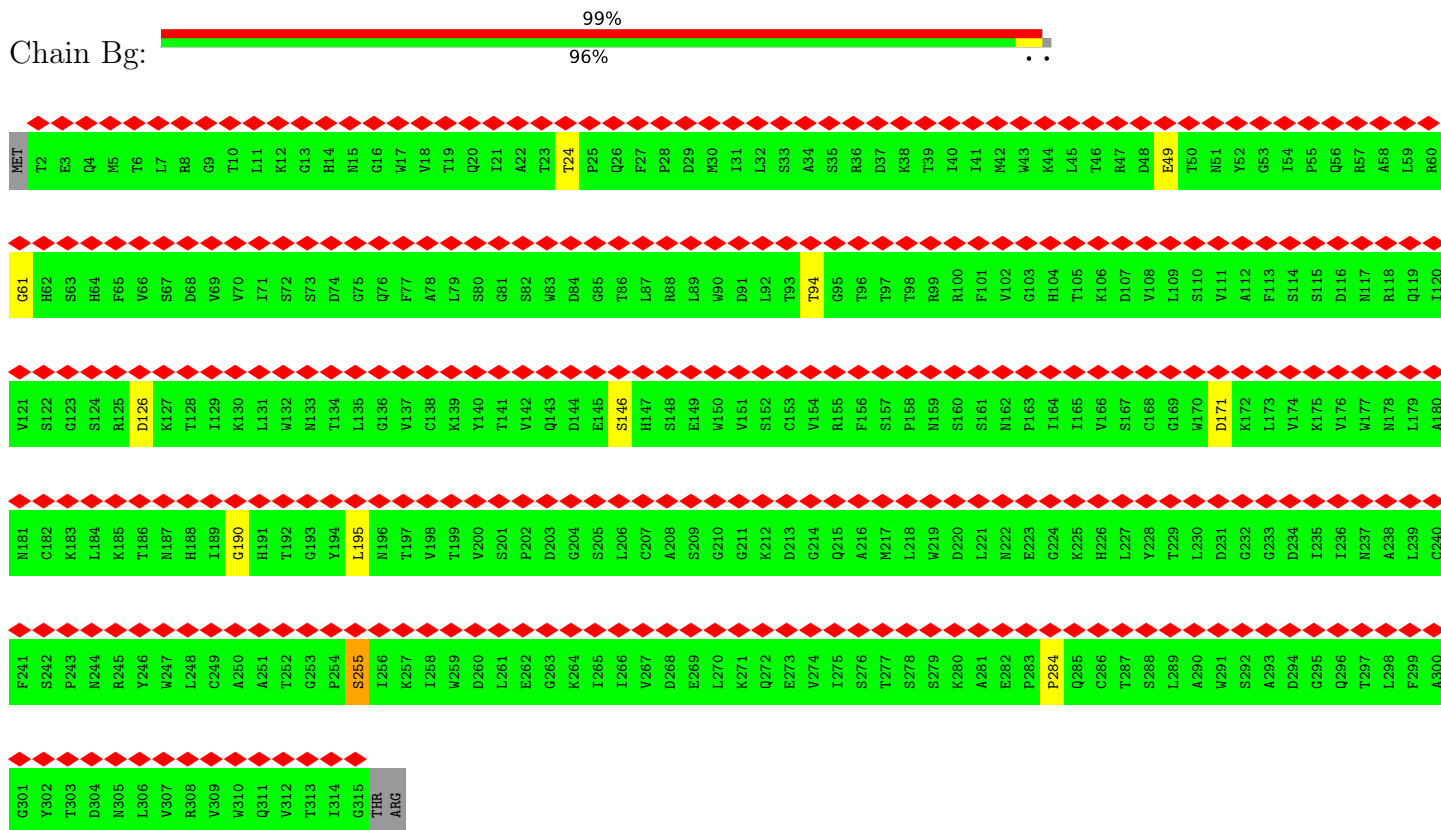


- Molecule 81: Ubiquitin-40S ribosomal protein S27a

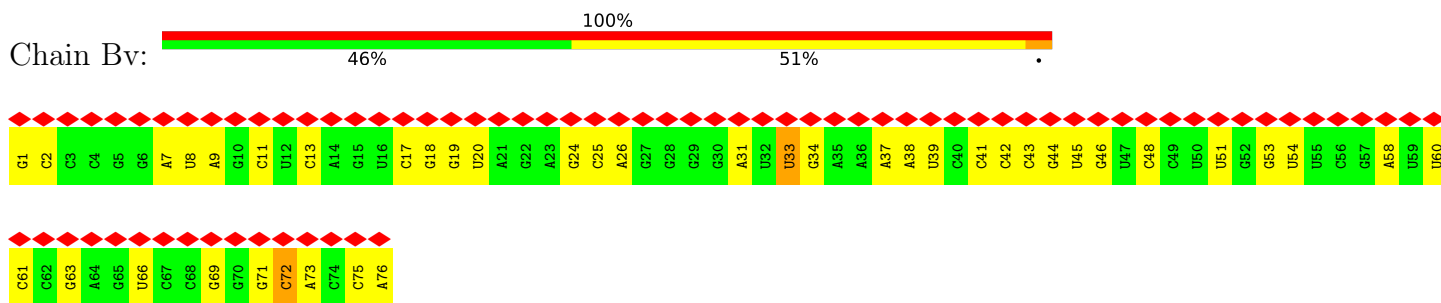




• Molecule 82: Guanine nucleotide-binding protein subunit beta-2-like 1



• Molecule 83: tRNA

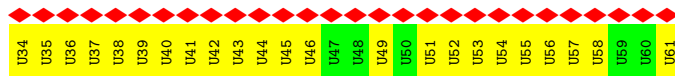


• Molecule 83: tRNA

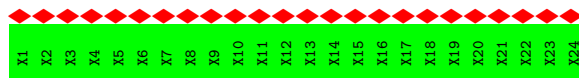




- Molecule 84: mRNA



- Molecule 85: Nascent protein chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	313321	Depositor
Resolution determination method	Not provided	
CTF correction method	DEFOCUS GROUPS	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20.00	Depositor
Minimum defocus (nm)	2000.00	Depositor
Maximum defocus (nm)	4500.00	Depositor
Magnification	205000	Depositor
Image detector	TVIPS TEMCAM-F416 (4k x 4k)	Depositor
Maximum map value	15.028	Depositor
Minimum map value	-7.978	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	3.0	Depositor
Map size (\AA)	378.0, 378.0, 378.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.945, 0.945, 0.945	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	A3	1.12	40/3726 (1.1%)	1.17	52/5804 (0.9%)
2	A4	1.10	32/2839 (1.1%)	1.13	40/4425 (0.9%)
3	AA	0.44	0/1968	0.65	1/2639 (0.0%)
4	AB	0.39	0/3246	0.64	0/4345
5	AC	0.41	1/2942 (0.0%)	0.62	3/3951 (0.1%)
6	AD	0.42	0/2437	0.65	2/3262 (0.1%)
7	AE	0.50	0/1603	0.80	4/2153 (0.2%)
8	AF	0.37	0/1986	0.57	0/2644
9	AG	0.38	0/1913	0.57	0/2576
10	AH	0.38	0/1545	0.54	1/2077 (0.0%)
11	AI	0.41	0/1730	0.61	1/2311 (0.0%)
12	AJ	0.39	0/1376	0.59	0/1841
13	AK	0.85	3/886 (0.3%)	1.53	23/1188 (1.9%)
14	AL	0.48	3/1688 (0.2%)	0.73	5/2260 (0.2%)
15	AM	0.39	0/1161	0.64	0/1554
16	AN	0.39	0/1746	0.57	0/2338
17	AO	0.37	0/1638	0.60	0/2191
18	AP	0.42	0/1268	0.70	0/1701
19	AQ	0.47	2/1537 (0.1%)	0.63	3/2052 (0.1%)
20	AR	0.38	0/1533	0.63	1/2025 (0.0%)
21	AS	0.38	1/1488 (0.1%)	0.60	1/1997 (0.1%)
22	AT	0.38	0/1312	0.60	0/1753
23	AU	0.35	0/822	0.58	0/1103
24	AV	0.37	0/983	0.55	0/1319
25	AW	0.42	1/1004 (0.1%)	0.74	3/1332 (0.2%)
26	AX	0.34	0/975	0.51	0/1312
27	AY	0.35	0/1081	0.56	0/1439
28	AZ	0.42	0/1126	0.69	2/1502 (0.1%)
29	Aa	0.50	0/1191	0.70	1/1591 (0.1%)
30	Ab	0.36	0/569	0.58	0/750
31	Ac	0.38	0/812	0.60	0/1089
32	Ad	0.38	0/894	0.58	0/1204

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	Ae	0.42	0/1082	0.58	1/1443 (0.1%)
34	Af	0.47	0/895	0.73	2/1198 (0.2%)
35	Ag	0.39	0/916	0.58	0/1220
36	Ah	0.37	0/1023	0.63	1/1351 (0.1%)
37	Ai	0.35	0/805	0.58	0/1065
38	Aj	0.47	1/703 (0.1%)	0.92	3/929 (0.3%)
39	Ak	0.51	0/575	0.74	0/761
40	Al	0.37	0/454	0.54	0/599
41	Am	0.42	0/417	0.59	0/553
42	An	0.36	0/241	0.60	0/305
43	Ao	0.39	0/877	0.66	0/1156
44	Ap	0.39	0/718	0.60	0/953
45	Aq	0.89	3/1058 (0.3%)	1.93	34/1424 (2.4%)
46	At	0.41	0/995	0.69	0/1334
47	Au	0.73	1/1772 (0.1%)	1.28	17/2375 (0.7%)
48	A2	0.98	666/86613 (0.8%)	1.13	1094/135108 (0.8%)
49	B1	0.96	275/40767 (0.7%)	1.13	515/63536 (0.8%)
50	BA	0.52	1/1741 (0.1%)	0.72	5/2366 (0.2%)
51	BB	0.42	0/1749	0.62	0/2340
52	BC	0.37	0/1761	0.58	0/2379
53	BD	0.44	1/1736 (0.1%)	0.63	1/2338 (0.0%)
54	BE	0.41	1/2072 (0.0%)	0.59	1/2793 (0.0%)
55	BF	0.40	1/1524 (0.1%)	0.62	1/2048 (0.0%)
56	BG	0.41	0/1907	0.62	1/2538 (0.0%)
57	BH	0.43	0/1501	0.64	1/2009 (0.0%)
58	BI	0.41	0/1725	0.59	0/2298
59	BJ	0.36	0/1520	0.55	1/2030 (0.0%)
60	BK	0.43	0/851	0.68	0/1147
61	BL	0.44	1/1281 (0.1%)	0.65	2/1710 (0.1%)
62	BM	0.39	0/941	0.63	0/1264
63	BN	0.40	1/1226 (0.1%)	0.57	1/1649 (0.1%)
64	BO	0.43	0/1029	0.64	0/1380
65	BP	0.50	1/1019 (0.1%)	0.74	3/1361 (0.2%)
66	BQ	0.38	0/1126	0.58	2/1506 (0.1%)
67	BR	0.49	3/1023 (0.3%)	0.68	3/1373 (0.2%)
68	BS	0.42	1/1172 (0.1%)	0.63	1/1570 (0.1%)
69	BT	0.36	0/1131	0.60	0/1515
70	BU	0.45	0/778	0.65	1/1045 (0.1%)
71	BV	0.41	0/623	0.60	1/833 (0.1%)
72	BW	0.37	0/1051	0.55	0/1406
73	BX	0.44	1/1097 (0.1%)	0.61	1/1464 (0.1%)
74	BY	0.40	0/1032	0.64	0/1371
75	BZ	0.43	0/696	0.62	0/929

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
76	Ba	0.49	0/786	0.75	2/1053 (0.2%)
77	Bb	0.42	0/637	0.59	0/854
78	Bc	0.40	0/490	0.66	0/656
79	Bd	0.54	0/437	0.89	3/580 (0.5%)
80	Be	0.37	0/443	0.60	0/583
81	Bf	0.44	0/613	0.62	0/811
82	Bg	0.40	0/2497	0.60	0/3399
83	Bv	0.58	4/1813 (0.2%)	0.84	5/2823 (0.2%)
83	Bw	0.43	2/1813 (0.1%)	0.90	4/2823 (0.1%)
84	Bx	0.60	2/616 (0.3%)	1.20	11/948 (1.2%)
All	All	0.80	1049/234393 (0.4%)	0.98	1861/344230 (0.5%)

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
4	AB	0	3
5	AC	0	2
7	AE	0	1
8	AF	0	2
10	AH	0	3
13	AK	0	3
14	AL	0	6
39	Ak	0	1
48	A2	2	3
49	B1	1	0
54	BE	0	6
63	BN	0	2
64	BO	0	2
65	BP	0	11
66	BQ	0	6
72	BW	0	3
76	Ba	0	1
79	Bd	0	1
All	All	3	56

The worst 5 of 1049 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
48	A2	3612	U	O3'-P	-21.32	1.35	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
49	B1	842	C	O3'-P	20.18	1.85	1.61
48	A2	2689	C	O3'-P	-19.59	1.37	1.61
49	B1	558	G	O3'-P	19.39	1.84	1.61
49	B1	497	C	O3'-P	18.59	1.83	1.61

The worst 5 of 1861 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
45	Aq	75	PRO	CA-N-CD	-36.28	60.70	111.50
48	A2	1225	G	O5'-P-OP1	-28.99	75.91	110.70
48	A2	131	C	C4'-C3'-O3'	28.78	170.56	113.00
49	B1	558	G	P-O3'-C3'	-28.42	85.60	119.70
48	A2	137	G	C4'-C3'-O3'	26.49	165.98	113.00

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
48	A2	131	C	C3'
48	A2	137	G	C3'
49	B1	1289	U	C4'

5 of 56 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	AB	295	ASP	Peptide
4	AB	296	GLY	Peptide
4	AB	297	LYS	Peptide
5	AC	98	GLY	Peptide
5	AC	99	GLY	Peptide

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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 PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	AA	250/257 (97%)	237 (95%)	7 (3%)	6 (2%)	6	35
4	AB	392/403 (97%)	370 (94%)	10 (3%)	12 (3%)	4	30
5	AC	361/427 (84%)	333 (92%)	13 (4%)	15 (4%)	3	23
6	AD	292/297 (98%)	270 (92%)	14 (5%)	8 (3%)	5	33
7	AE	192/288 (67%)	161 (84%)	12 (6%)	19 (10%)	0	7
8	AF	232/248 (94%)	213 (92%)	10 (4%)	9 (4%)	3	25
9	AG	232/266 (87%)	217 (94%)	10 (4%)	5 (2%)	6	37
10	AH	189/192 (98%)	175 (93%)	8 (4%)	6 (3%)	4	29
11	AI	204/214 (95%)	192 (94%)	7 (3%)	5 (2%)	5	34
12	AJ	167/178 (94%)	154 (92%)	7 (4%)	6 (4%)	3	26
13	AK	107/317 (34%)	34 (32%)	37 (35%)	36 (34%)	0	0
14	AL	203/211 (96%)	175 (86%)	14 (7%)	14 (7%)	1	12
15	AM	137/215 (64%)	127 (93%)	5 (4%)	5 (4%)	3	26
16	AN	201/204 (98%)	193 (96%)	6 (3%)	2 (1%)	15	54
17	AO	193/203 (95%)	187 (97%)	3 (2%)	3 (2%)	9	43
18	AP	151/184 (82%)	147 (97%)	4 (3%)	0	100	100
19	AQ	185/188 (98%)	163 (88%)	10 (5%)	12 (6%)	1	14
20	AR	179/196 (91%)	171 (96%)	4 (2%)	4 (2%)	6	37
21	AS	173/176 (98%)	156 (90%)	11 (6%)	6 (4%)	3	27
22	AT	155/160 (97%)	144 (93%)	6 (4%)	5 (3%)	4	29
23	AU	97/128 (76%)	81 (84%)	7 (7%)	9 (9%)	0	8
24	AV	127/140 (91%)	125 (98%)	2 (2%)	0	100	100
25	AW	119/157 (76%)	96 (81%)	18 (15%)	5 (4%)	3	23
26	AX	115/156 (74%)	113 (98%)	2 (2%)	0	100	100
27	AY	125/145 (86%)	119 (95%)	2 (2%)	4 (3%)	4	29
28	AZ	132/136 (97%)	118 (89%)	6 (4%)	8 (6%)	1	15
29	Aa	145/148 (98%)	135 (93%)	6 (4%)	4 (3%)	5	32
30	Ab	66/159 (42%)	57 (86%)	5 (8%)	4 (6%)	1	15
31	Ac	101/115 (88%)	97 (96%)	2 (2%)	2 (2%)	7	39
32	Ad	104/125 (83%)	99 (95%)	3 (3%)	2 (2%)	8	40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
33	Ae	127/135 (94%)	116 (91%)	6 (5%)	5 (4%)	3	25
34	Af	107/110 (97%)	93 (87%)	5 (5%)	9 (8%)	1	9
35	Ag	112/117 (96%)	107 (96%)	3 (3%)	2 (2%)	8	41
36	Ah	120/123 (98%)	112 (93%)	2 (2%)	6 (5%)	2	19
37	Ai	95/105 (90%)	83 (87%)	7 (7%)	5 (5%)	2	17
38	Aj	82/97 (84%)	69 (84%)	7 (8%)	6 (7%)	1	11
39	Ak	67/70 (96%)	50 (75%)	7 (10%)	10 (15%)	0	3
40	Al	48/51 (94%)	47 (98%)	1 (2%)	0	100	100
41	Am	48/128 (38%)	44 (92%)	3 (6%)	1 (2%)	7	38
42	An	23/25 (92%)	23 (100%)	0	0	100	100
43	Ao	103/106 (97%)	94 (91%)	5 (5%)	4 (4%)	3	25
44	Ap	89/92 (97%)	84 (94%)	3 (3%)	2 (2%)	6	37
45	Aq	136/165 (82%)	40 (29%)	47 (35%)	49 (36%)	0	0
46	At	120/137 (88%)	107 (89%)	10 (8%)	3 (2%)	5	34
47	Au	215/217 (99%)	185 (86%)	20 (9%)	10 (5%)	2	20
50	BA	213/295 (72%)	197 (92%)	12 (6%)	4 (2%)	8	40
51	BB	210/264 (80%)	180 (86%)	12 (6%)	18 (9%)	1	9
52	BC	220/293 (75%)	204 (93%)	7 (3%)	9 (4%)	3	23
53	BD	218/243 (90%)	201 (92%)	10 (5%)	7 (3%)	4	29
54	BE	255/263 (97%)	230 (90%)	14 (6%)	11 (4%)	2	22
55	BF	188/204 (92%)	163 (87%)	15 (8%)	10 (5%)	2	17
56	BG	230/249 (92%)	211 (92%)	11 (5%)	8 (4%)	3	27
57	BH	181/194 (93%)	169 (93%)	8 (4%)	4 (2%)	6	37
58	BI	205/208 (99%)	175 (85%)	20 (10%)	10 (5%)	2	19
59	BJ	177/194 (91%)	137 (77%)	28 (16%)	12 (7%)	1	13
60	BK	96/165 (58%)	84 (88%)	7 (7%)	5 (5%)	2	18
61	BL	151/158 (96%)	133 (88%)	11 (7%)	7 (5%)	2	21
62	BM	118/132 (89%)	113 (96%)	1 (1%)	4 (3%)	3	28
63	BN	147/151 (97%)	126 (86%)	13 (9%)	8 (5%)	2	17
64	BO	134/151 (89%)	113 (84%)	11 (8%)	10 (8%)	1	11
65	BP	118/145 (81%)	100 (85%)	9 (8%)	9 (8%)	1	10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
66	BQ	137/146 (94%)	120 (88%)	10 (7%)	7 (5%)	2	19
67	BR	123/135 (91%)	106 (86%)	9 (7%)	8 (6%)	1	14
68	BS	137/152 (90%)	125 (91%)	7 (5%)	5 (4%)	3	26
69	BT	141/145 (97%)	131 (93%)	6 (4%)	4 (3%)	5	32
70	BU	95/119 (80%)	91 (96%)	2 (2%)	2 (2%)	7	38
71	BV	79/83 (95%)	77 (98%)	2 (2%)	0	100	100
72	BW	127/130 (98%)	120 (94%)	3 (2%)	4 (3%)	4	30
73	BX	137/143 (96%)	124 (90%)	8 (6%)	5 (4%)	3	26
74	BY	123/133 (92%)	116 (94%)	6 (5%)	1 (1%)	19	58
75	BZ	84/125 (67%)	80 (95%)	1 (1%)	3 (4%)	3	26
76	Ba	95/115 (83%)	87 (92%)	7 (7%)	1 (1%)	14	52
77	Bb	78/84 (93%)	72 (92%)	4 (5%)	2 (3%)	5	33
78	Bc	60/69 (87%)	57 (95%)	1 (2%)	2 (3%)	4	28
79	Bd	49/56 (88%)	42 (86%)	4 (8%)	3 (6%)	1	15
80	Be	53/59 (90%)	49 (92%)	2 (4%)	2 (4%)	3	25
81	Bf	71/156 (46%)	63 (89%)	7 (10%)	1 (1%)	11	46
82	Bg	312/317 (98%)	292 (94%)	12 (4%)	8 (3%)	5	33
All	All	11580/13387 (86%)	10401 (90%)	657 (6%)	522 (4%)	4	21

5 of 522 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AA	138	SER
3	AA	144	LYS
3	AA	197	PRO
4	AB	189	THR
4	AB	356	LYS

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	AA	194/199 (98%)	185 (95%)	9 (5%)	27	61
4	AB	343/349 (98%)	323 (94%)	20 (6%)	20	53
5	AC	302/348 (87%)	278 (92%)	24 (8%)	12	41
6	AD	248/250 (99%)	229 (92%)	19 (8%)	13	42
7	AE	174/252 (69%)	127 (73%)	47 (27%)	0	3
8	AF	203/215 (94%)	193 (95%)	10 (5%)	25	59
9	AG	199/223 (89%)	192 (96%)	7 (4%)	36	67
10	AH	170/171 (99%)	160 (94%)	10 (6%)	19	53
11	AI	178/181 (98%)	174 (98%)	4 (2%)	52	78
12	AJ	142/149 (95%)	138 (97%)	4 (3%)	43	72
13	AK	95/258 (37%)	70 (74%)	25 (26%)	0	3
14	AL	171/177 (97%)	156 (91%)	15 (9%)	10	38
15	AM	118/161 (73%)	104 (88%)	14 (12%)	5	25
16	AN	171/172 (99%)	163 (95%)	8 (5%)	26	60
17	AO	168/174 (97%)	164 (98%)	4 (2%)	49	76
18	AP	134/163 (82%)	123 (92%)	11 (8%)	11	40
19	AQ	164/165 (99%)	144 (88%)	20 (12%)	5	23
20	AR	160/175 (91%)	145 (91%)	15 (9%)	8	35
21	AS	156/157 (99%)	149 (96%)	7 (4%)	27	61
22	AT	138/140 (99%)	126 (91%)	12 (9%)	10	38
23	AU	89/115 (77%)	78 (88%)	11 (12%)	4	23
24	AV	100/107 (94%)	98 (98%)	2 (2%)	55	79
25	AW	100/126 (79%)	94 (94%)	6 (6%)	19	52
26	AX	105/133 (79%)	95 (90%)	10 (10%)	8	34
27	AY	119/135 (88%)	109 (92%)	10 (8%)	11	40
28	AZ	117/118 (99%)	100 (86%)	17 (14%)	3	18
29	Aa	120/121 (99%)	110 (92%)	10 (8%)	11	40
30	Ab	58/126 (46%)	54 (93%)	4 (7%)	15	47
31	Ac	88/97 (91%)	85 (97%)	3 (3%)	37	68
32	Ad	97/110 (88%)	96 (99%)	1 (1%)	76	88
33	Ae	115/121 (95%)	114 (99%)	1 (1%)	78	90
34	Af	88/89 (99%)	85 (97%)	3 (3%)	37	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
35	Ag	98/100 (98%)	93 (95%)	5 (5%)	24	57
36	Ah	109/110 (99%)	96 (88%)	13 (12%)	5	25
37	Ai	83/89 (93%)	80 (96%)	3 (4%)	35	66
38	Aj	71/80 (89%)	65 (92%)	6 (8%)	10	39
39	Ak	64/65 (98%)	46 (72%)	18 (28%)	0	3
40	Al	47/48 (98%)	44 (94%)	3 (6%)	17	50
41	Am	46/116 (40%)	45 (98%)	1 (2%)	52	78
42	An	24/24 (100%)	23 (96%)	1 (4%)	30	63
43	Ao	93/94 (99%)	85 (91%)	8 (9%)	10	38
44	Ap	74/75 (99%)	71 (96%)	3 (4%)	30	63
45	Aq	114/137 (83%)	90 (79%)	24 (21%)	1	5
46	At	106/121 (88%)	88 (83%)	18 (17%)	2	12
47	Au	196/196 (100%)	180 (92%)	16 (8%)	11	40
50	BA	180/243 (74%)	165 (92%)	15 (8%)	11	40
51	BB	193/231 (84%)	177 (92%)	16 (8%)	11	40
52	BC	188/225 (84%)	177 (94%)	11 (6%)	19	53
53	BD	183/202 (91%)	175 (96%)	8 (4%)	28	62
54	BE	220/225 (98%)	213 (97%)	7 (3%)	39	69
55	BF	160/170 (94%)	155 (97%)	5 (3%)	40	70
56	BG	202/218 (93%)	177 (88%)	25 (12%)	4	23
57	BH	164/174 (94%)	160 (98%)	4 (2%)	49	76
58	BI	179/180 (99%)	175 (98%)	4 (2%)	52	78
59	BJ	160/168 (95%)	151 (94%)	9 (6%)	21	54
60	BK	89/136 (65%)	83 (93%)	6 (7%)	16	48
61	BL	138/142 (97%)	134 (97%)	4 (3%)	42	71
62	BM	102/108 (94%)	101 (99%)	1 (1%)	76	88
63	BN	130/131 (99%)	122 (94%)	8 (6%)	18	51
64	BO	106/119 (89%)	99 (93%)	7 (7%)	16	49
65	BP	109/130 (84%)	93 (85%)	16 (15%)	3	18
66	BQ	115/121 (95%)	102 (89%)	13 (11%)	6	27
67	BR	113/122 (93%)	97 (86%)	16 (14%)	3	19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
68	BS	121/132 (92%)	106 (88%)	15 (12%)	4	23
69	BT	113/115 (98%)	102 (90%)	11 (10%)	8	33
70	BU	90/107 (84%)	85 (94%)	5 (6%)	21	54
71	BV	65/67 (97%)	59 (91%)	6 (9%)	9	36
72	BW	112/113 (99%)	108 (96%)	4 (4%)	35	66
73	BX	111/115 (96%)	101 (91%)	10 (9%)	9	37
74	BY	107/115 (93%)	103 (96%)	4 (4%)	34	65
75	BZ	75/103 (73%)	57 (76%)	18 (24%)	0	4
76	Ba	84/98 (86%)	75 (89%)	9 (11%)	6	30
77	Bb	72/76 (95%)	72 (100%)	0	100	100
78	Bc	55/62 (89%)	52 (94%)	3 (6%)	21	54
79	Bd	45/49 (92%)	40 (89%)	5 (11%)	6	28
80	Be	44/48 (92%)	38 (86%)	6 (14%)	3	20
81	Bf	66/140 (47%)	65 (98%)	1 (2%)	65	84
82	Bg	272/275 (99%)	268 (98%)	4 (2%)	65	84
All	All	10112/11392 (89%)	9354 (92%)	758 (8%)	17	43

5 of 758 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
46	At	89	THR
58	BI	19	LYS
47	Au	185	LEU
46	At	87	ARG
52	BC	166	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 196 such sidechains are listed below:

Mol	Chain	Res	Type
46	At	70	GLN
57	BH	114	GLN
47	Au	143	ASN
52	BC	235	ASN
61	BL	13	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A3	156/194 (80%)	28 (17%)	4 (2%)
2	A4	118/121 (97%)	23 (19%)	1 (0%)
48	A2	3600/5029 (71%)	695 (19%)	48 (1%)
49	B1	1701/1869 (91%)	290 (17%)	19 (1%)
83	Bv	75/76 (98%)	38 (50%)	0
83	Bw	75/76 (98%)	34 (45%)	0
84	Bx	27/28 (96%)	18 (66%)	0
All	All	5752/7393 (77%)	1126 (19%)	72 (1%)

5 of 1126 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A3	2	G
1	A3	12	G
1	A3	16	G
1	A3	35	C
1	A3	59	A

5 of 72 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
49	B1	283	G
49	B1	1756	C
49	B1	369	C
49	B1	797	C
48	A2	1402	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 331 ligands modelled in this entry, 331 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
49	B1	8
48	A2	5

The worst 5 of 13 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B1	842:C	O3'	843:C	P	1.85
1	B1	558:G	O3'	559:G	P	1.84
1	B1	497:C	O3'	498:C	P	1.83
1	B1	72:C	O3'	73:C	P	1.82
1	B1	1253:A	O3'	1254:C	P	1.82

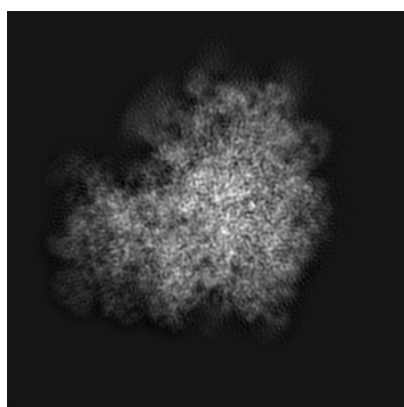
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-2875. These allow visual inspection of the internal detail of the map and identification of artifacts.

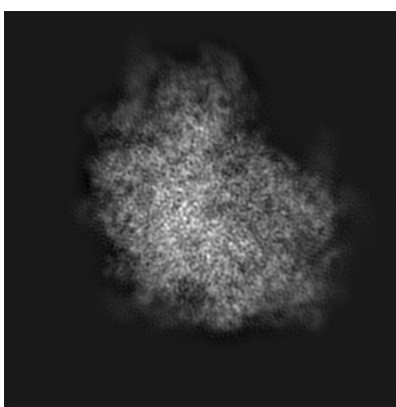
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

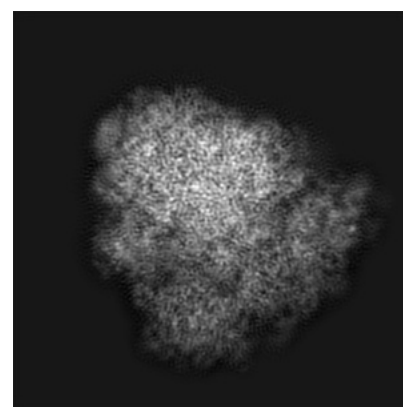
6.1.1 Primary map



X



Y

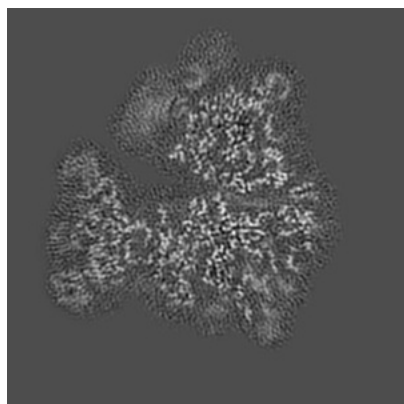


Z

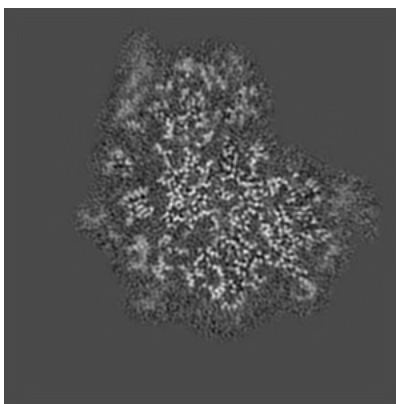
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

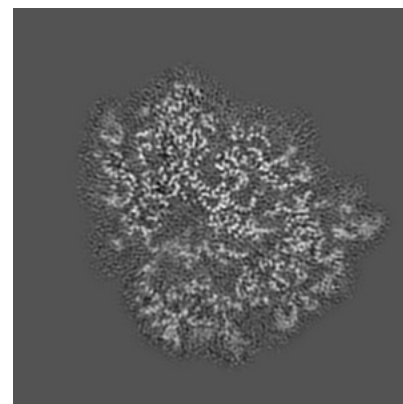
6.2.1 Primary map



X Index: 200



Y Index: 200

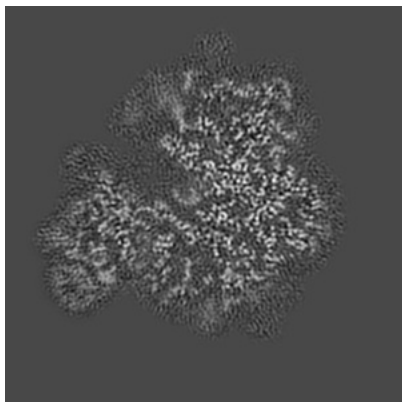


Z Index: 200

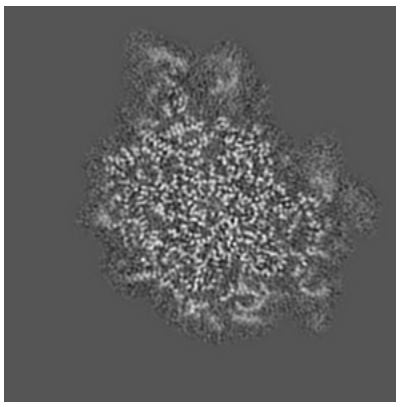
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

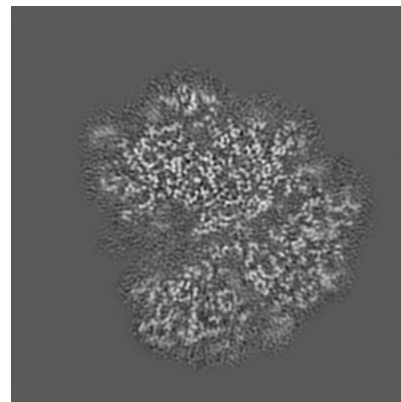
6.3.1 Primary map



X Index: 186



Y Index: 220

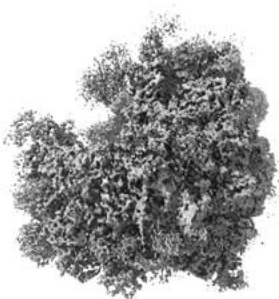


Z Index: 184

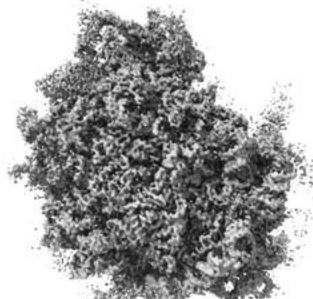
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

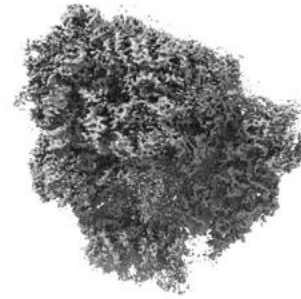
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 3.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

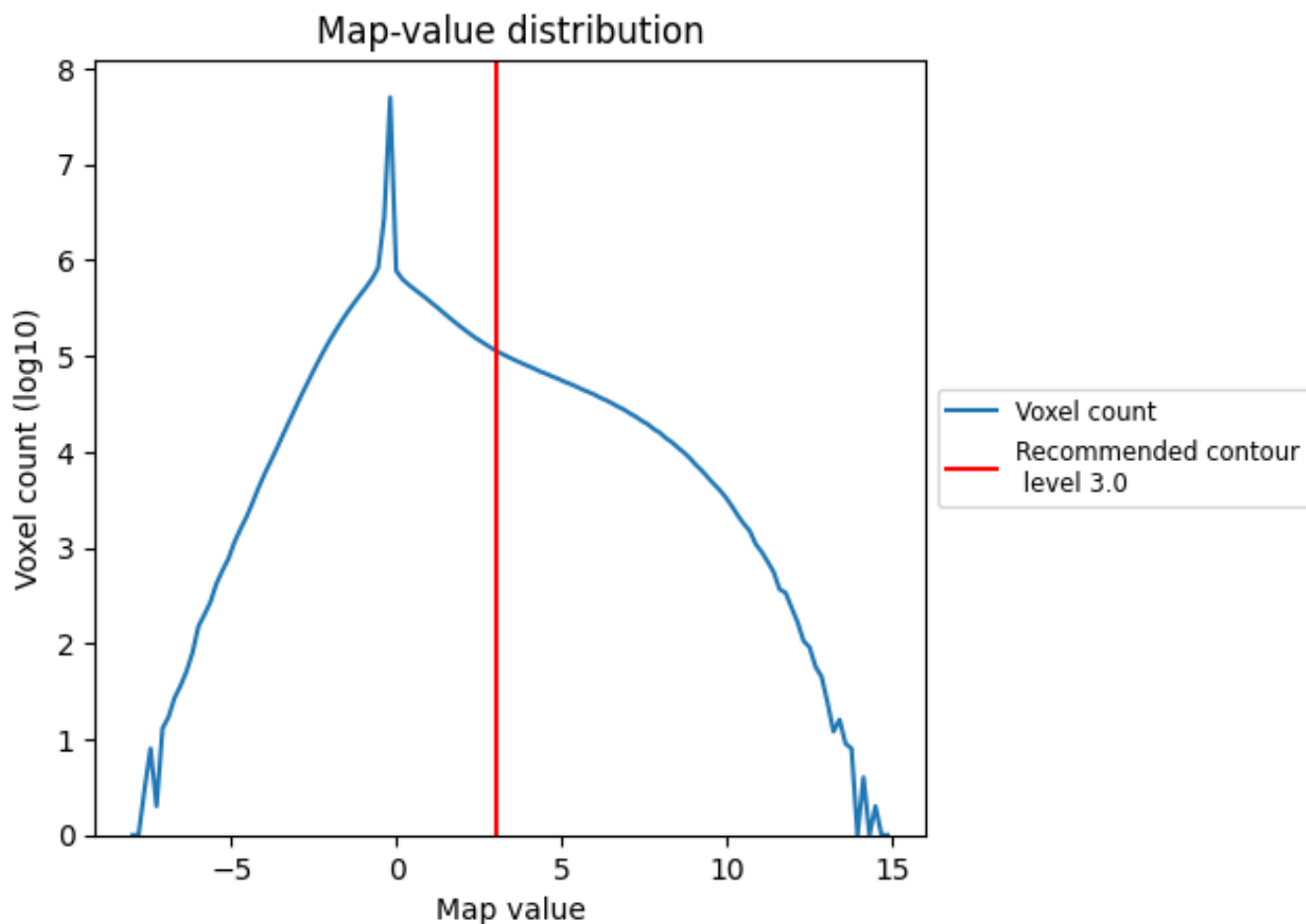
6.5 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

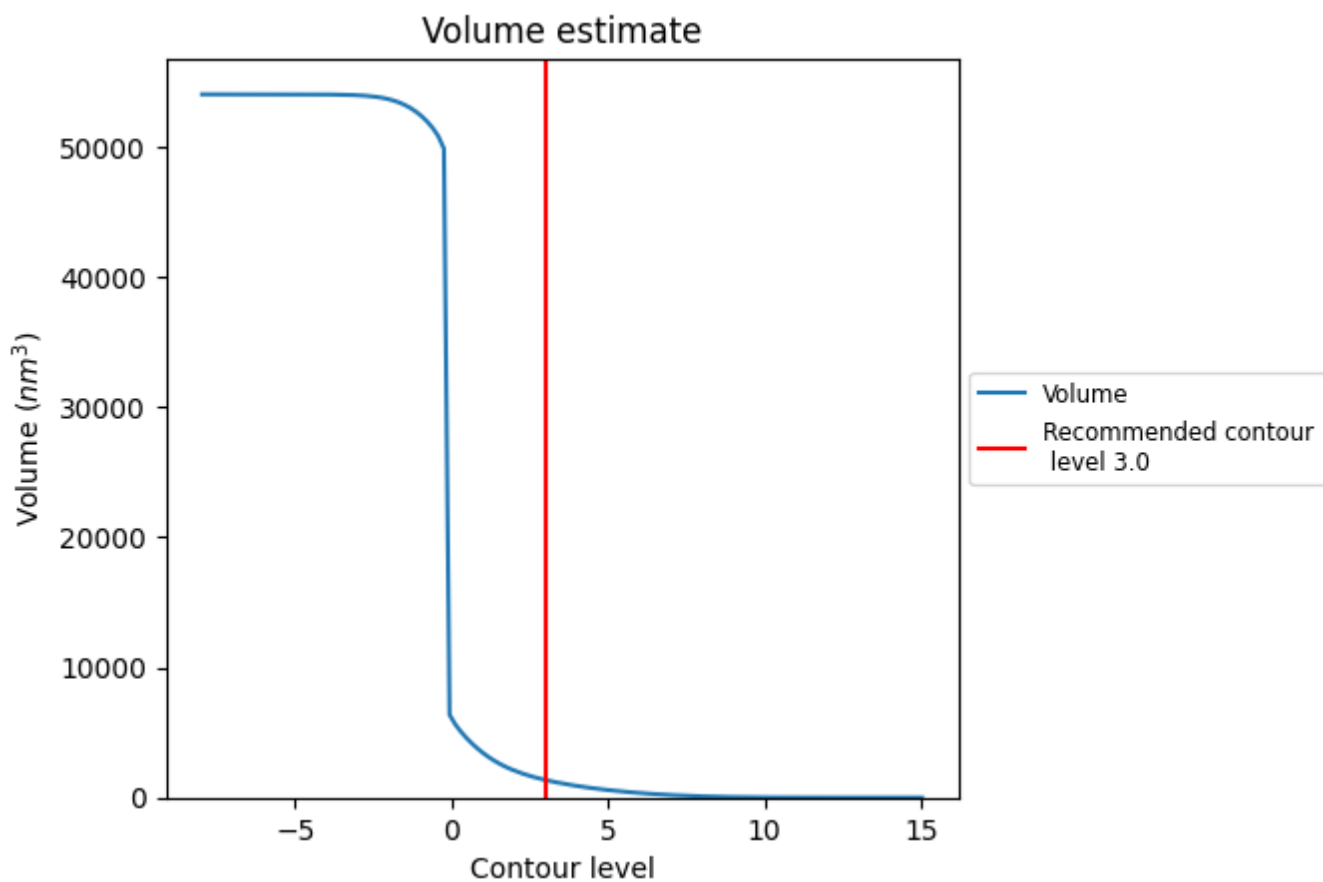
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

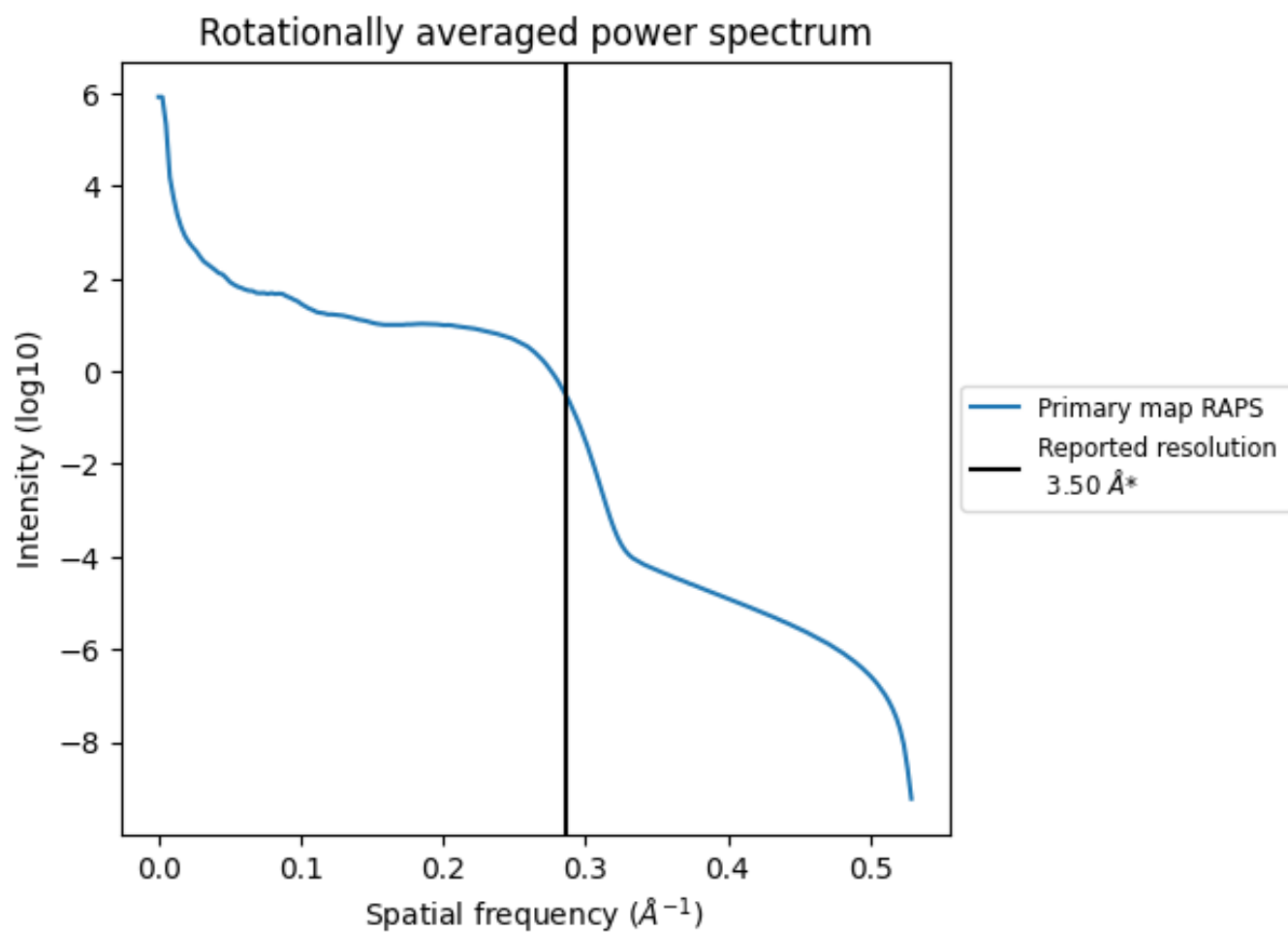
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 1346 nm³; this corresponds to an approximate mass of 1216 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum i



*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

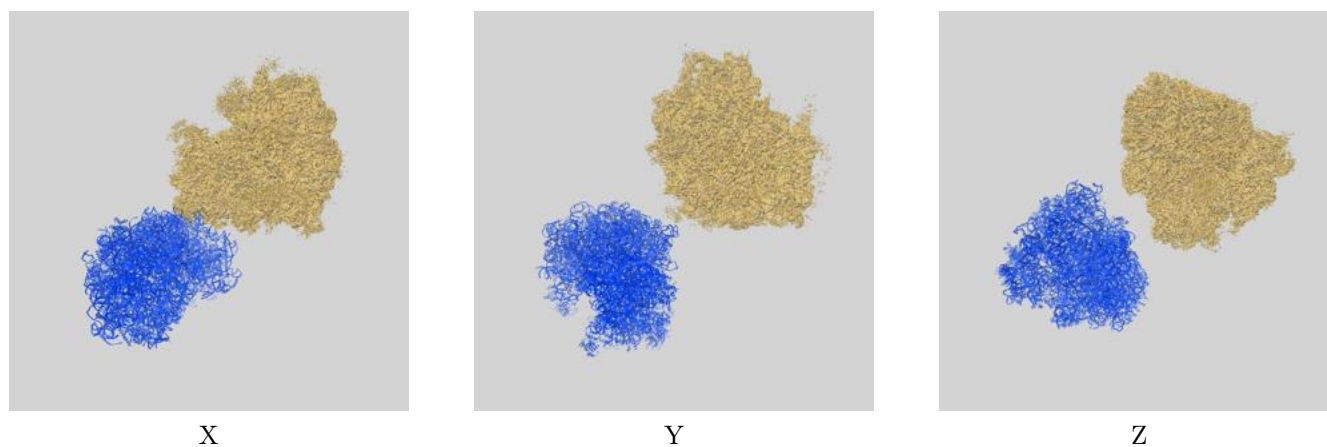
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

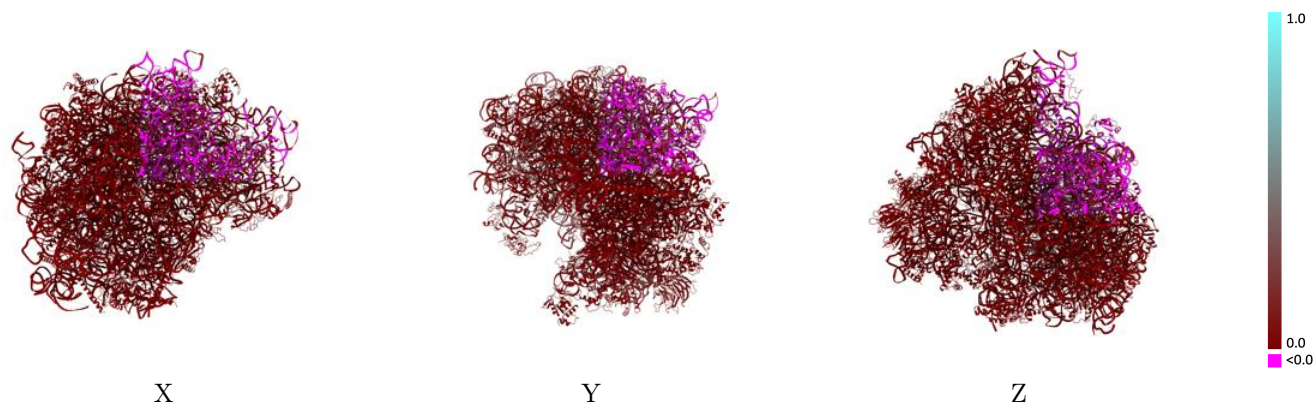
This section contains information regarding the fit between EMDB map EMD-2875 and PDB model 5AJ0. Per-residue inclusion information can be found in section 3 on page 21.

9.1 Map-model overlay [i](#)



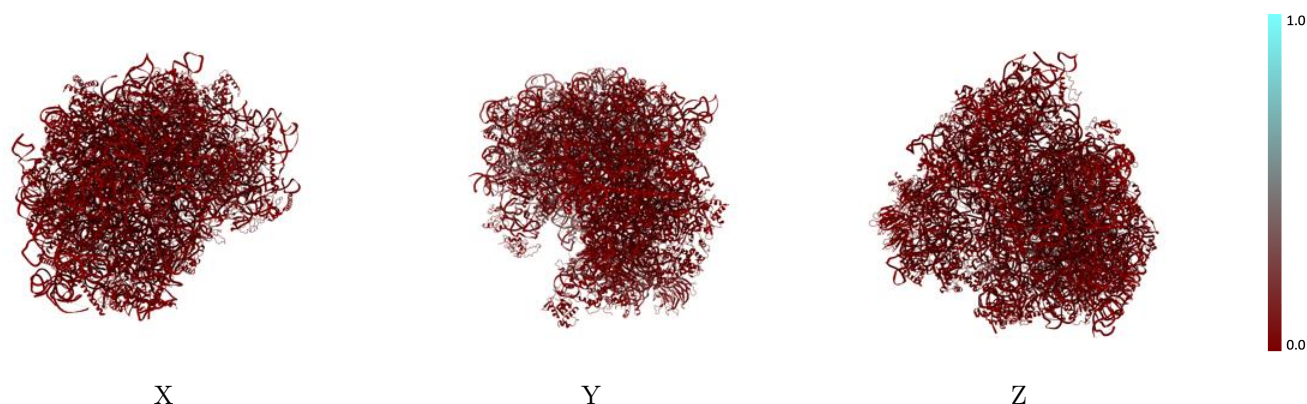
The images above show the 3D surface view of the map at the recommended contour level 3.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



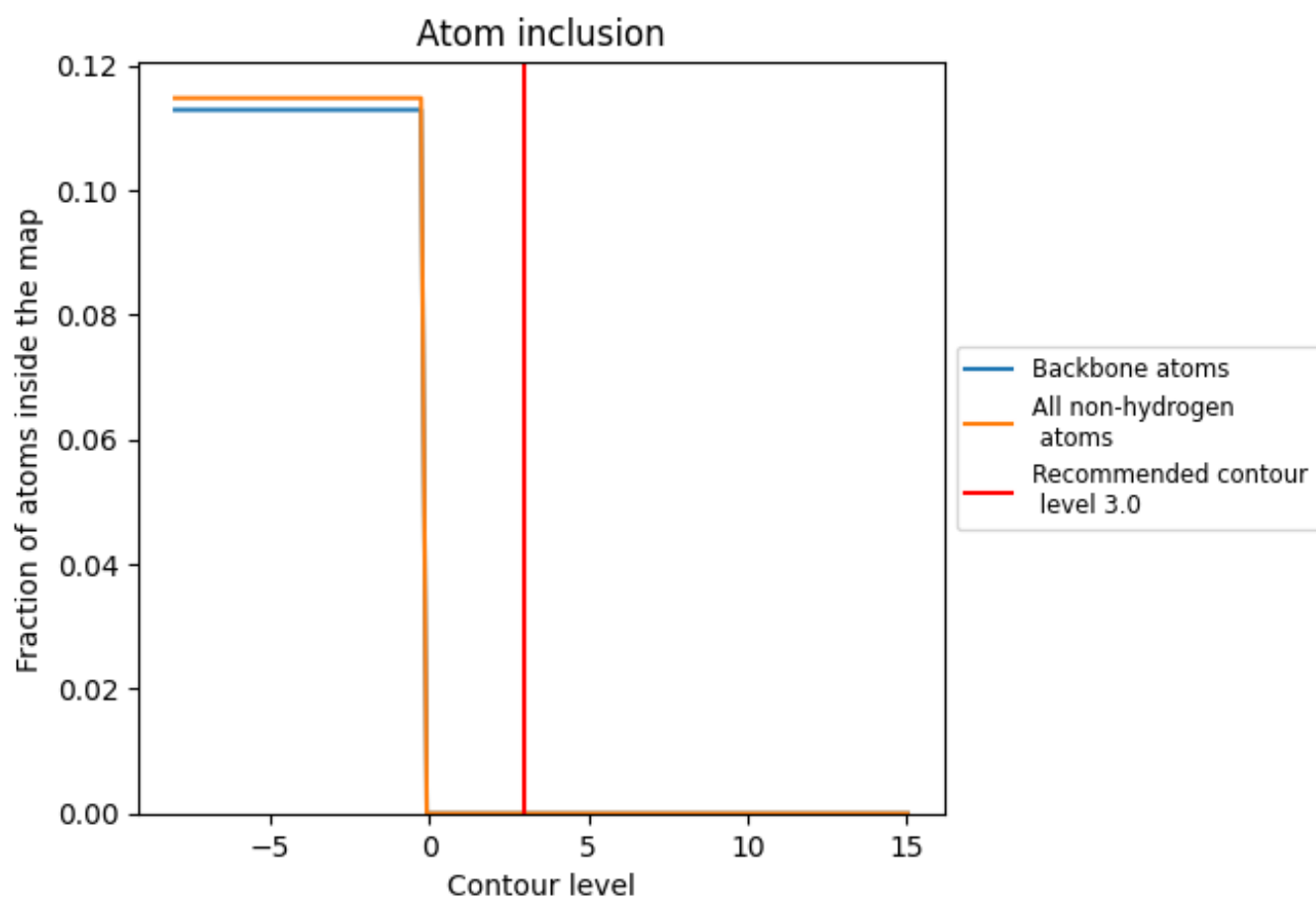
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (3.0).

9.4 Atom inclusion [i](#)



At the recommended contour level, 0% of all backbone atoms, 0% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (3.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.0000	0.0000
A2	0.0000	0.0000
A3	0.0000	0.0030
A4	0.0000	0.0000
AA	0.0000	0.0010
AB	0.0000	0.0000
AC	0.0000	0.0000
AD	0.0000	0.0000
AE	0.0000	0.0000
AF	0.0000	0.0000
AG	0.0000	-0.0010
AH	0.0000	0.0000
AI	0.0000	0.0000
AJ	0.0000	0.0000
AK	0.0000	0.0000
AL	0.0000	0.0000
AM	0.0000	0.0000
AN	0.0000	0.0000
AO	0.0000	0.0000
AP	0.0000	0.0010
AQ	0.0000	0.0000
AR	0.0000	0.0150
AS	0.0000	0.0000
AT	0.0000	0.0000
AU	0.0000	-0.0210
AV	0.0000	0.0000
AW	0.0000	0.0000
AX	0.0000	0.0020
AY	0.0000	0.0000
AZ	0.0000	0.0050
Aa	0.0000	0.0000
Ab	0.0000	0.0000
Ac	0.0000	-0.0080
Ad	0.0000	0.0000
Ae	0.0000	0.0000























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Chain	Atom inclusion	Q-score
Af	0.0000	0.0000
Ag	0.0000	0.0190
Ah	0.0000	0.0000
Ai	0.0000	0.0000
Aj	0.0000	-0.0050
Ak	0.0000	0.0040
Al	0.0000	0.0220
Am	0.0000	0.0000
An	0.0000	0.0000
Ao	0.0000	0.0000
Ap	0.0000	-0.0110
Aq	0.0000	0.0000
At	0.0000	0.0000
Au	0.0000	0.0000
B1	0.0000	0.0000
BA	0.0000	0.0000
BB	0.0000	0.0000
BC	0.0000	0.0000
BD	0.0000	0.0000
BE	0.0000	0.0000
BF	0.0000	0.0000
BG	0.0000	0.0000
BH	0.0000	0.0000
BI	0.0000	0.0150
BJ	0.0000	0.0000
BK	0.0000	0.0000
BL	0.0000	0.0040
BM	0.0000	0.0000
BN	0.0000	0.0010
BO	0.0000	0.0000
BP	0.0000	0.0000
BQ	0.0000	0.0000
BR	0.0000	0.0000
BS	0.0000	0.0000
BT	0.0000	0.0000
BU	0.0000	0.0000
BV	0.0000	0.0000
BW	0.0000	0.0000
BX	0.0000	0.0000
BY	0.0000	0.0000
BZ	0.0000	0.0000
Ba	0.0000	0.0000

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Chain	Atom inclusion	Q-score
Bb	 0.0000	 0.0000
Bc	 0.0000	 0.0000
Bd	 0.0000	 0.0000
Be	 0.0000	 0.0000
Bf	 0.0000	 0.0000
Bg	 0.0000	 0.0000
Bv	 0.0000	 0.0000
Bw	 0.0000	 0.0000
Bx	 0.0000	 0.0000
By	 0.0000	 0.0220