



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 10:42 AM UTC

PDB ID : 6LQQ / pdb_00006lqq
EMDB ID : EMD-0950
Title : Cryo-EM structure of 90S small subunit preribosomes in transition states (State B)
Authors : Du, Y.; Ye, K.
Deposited on : 2020-01-14
Resolution : 4.10 Å(reported)
Based on initial model : 6LQP

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:

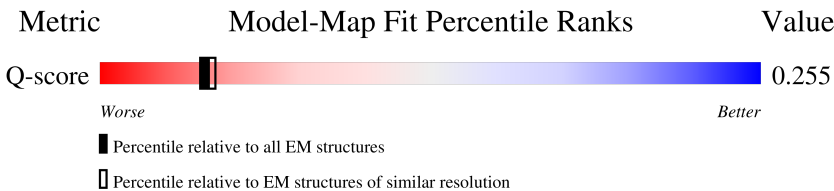
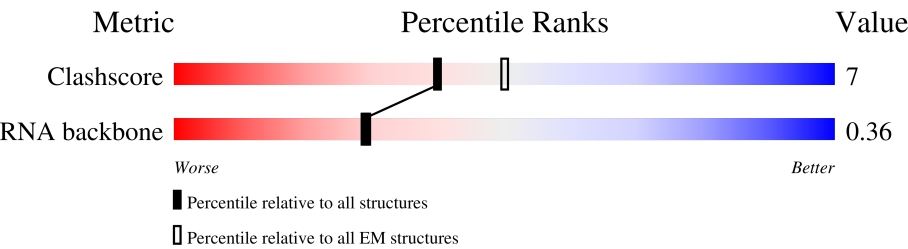
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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 4.10 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
RNA backbone	8273	3508	-
Q-score	-	25397	6458 (3.60 - 4.60)

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	3A	333	11% 34% 17% • 47%
2	5A	700	41% 33% 18% • 46%
3	SA	1808	27% 36% 31% 6% 27%
4	SC	255	23% 71% 19% 10%
5	SF	261	26% 64% 24% 12%

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Mol	Chain	Length	Quality of chain
6	SG	225	
7	SH	236	
8	SI	190	
9	SJ	200	
10	SK	197	
11	SM	156	
12	SO	151	
13	SP	137	
14	SR	143	
15	SX	130	
16	SY	145	
17	SZ	135	
18	Sc	82	
19	Sd	67	
20	3B	327	
20	3C	327	
21	3D	504	
22	3E	511	
23	3F	573	
24	3G	126	
24	3H	126	
25	A4	776	
26	A5	643	
27	A8	713	
28	A9	575	

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Mol	Chain	Length	Quality of chain
29	AE	1769	
30	AF	513	
31	AG	896	
32	B1	923	
33	B2	943	
34	B3	817	
35	B8	594	
36	BE	939	
37	B6	440	
38	5B	214	
39	5C	554	
40	5D	250	
41	5E	593	
42	5F	183	
43	5G	290	
44	5H	610	
45	5I	489	
46	5J	217	
47	5K	189	
48	RA	707	
49	RB	357	
50	RC	316	
51	RD	1729	
52	RE	1237	
53	RF	297	

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Mol	Chain	Length	Quality of chain
54	RG	252	
54	RH	252	
55	RJ	1183	
56	RK	367	
57	RL	1056	
57	RM	1056	
58	RN	810	
59	RO	552	
60	RP	2493	
61	RQ	899	
62	RS	483	
63	RT	326	
64	RV	346	
65	RY	534	
66	X1	347	

2 Entry composition

There are 69 unique types of molecules in this entry. The entry contains 226161 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 U3 snoRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	3A	175	Total	C	N	O	P	0	0
			3711	1661	648	1227	175		

- Molecule 2 is a RNA chain called 5' ETS.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	5A	375	Total	C	N	O	P	0	0
			8009	3578	1426	2630	375		

- Molecule 3 is a RNA chain called 18S pre-rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	SA	1323	Total	C	N	O	P	0	0
			28210	12608	5022	9257	1323		

- Molecule 4 is a protein called 40S ribosomal protein S1-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	SC	230	Total	C	N	O	S	0	0
			1830	1156	335	335	4		

- Molecule 5 is a protein called 40S ribosomal protein S4-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	SF	229	Total	C	N	O	S	0	0
			1815	1161	331	320	3		

- Molecule 6 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	SG	213	Total	C	N	O	S	0	0
			1669	1045	307	314	3		

- Molecule 7 is a protein called 40S ribosomal protein S6-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	SH	167	Total	C	N	O	S	0	0
			1327	834	256	235	2		

- Molecule 8 is a protein called 40S ribosomal protein S7-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	SI	165	Total	C	N	O		0	0
			1321	853	226	242			

- Molecule 9 is a protein called 40S ribosomal protein S8-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	SJ	166	Total	C	N	O	S	0	0
			1324	824	262	236	2		

- Molecule 10 is a protein called 40S ribosomal protein S9-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	SK	171	Total	C	N	O	S	0	0
			1388	879	268	240	1		

- Molecule 11 is a protein called 40S ribosomal protein S11-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	SM	123	Total	C	N	O	S	0	0
			997	641	189	164	3		

- Molecule 12 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	SO	134	Total	C	N	O	S	0	0
			1087	698	202	186	1		

- Molecule 13 is a protein called 40S ribosomal protein S14-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	SP	118	Total	C	N	O	S	0	0
			868	536	164	165	3		

- Molecule 14 is a protein called 40S ribosomal protein S16-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
14	SR	125	Total	C	N	O	0	0
			973	625	174	174		

- Molecule 15 is a protein called 40S ribosomal protein S22-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	SX	127	Total	C	N	O	S	0	0
			1003	640	183	177	3		

- Molecule 16 is a protein called 40S ribosomal protein S23-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	SY	103	Total	C	N	O	S	0	0
			786	503	144	137	2		

- Molecule 17 is a protein called 40S ribosomal protein S24-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
17	SZ	102	Total	C	N	O	0	0
			809	517	148	144		

- Molecule 18 is a protein called 40S ribosomal protein S27-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Sc	80	Total	C	N	O	S	0	0
			603	377	109	112	5		

- Molecule 19 is a protein called 40S ribosomal protein S28-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	Sd	63	Total	C	N	O	S	0	0
			497	306	99	91	1		

- Molecule 20 is a protein called rRNA 2'-O-methyltransferase fibrillarin.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	3B	240	Total	C	N	O	S	0	0
			1865	1184	333	338	10		
20	3C	225	Total	C	N	O	S	0	0
			1763	1120	316	317	10		

- Molecule 21 is a protein called Nucleolar protein 56.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	3D	369	Total	C	N	O	S	0	0
			2848	1811	489	540	8		

- Molecule 22 is a protein called Nucleolar protein 58.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	3E	431	Total	C	N	O	S	0	0
			3028	1888	543	588	9		

- Molecule 23 is a protein called Ribosomal RNA-processing protein 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	3F	454	Total	C	N	O	S	0	0
			3643	2315	638	680	10		

- Molecule 24 is a protein called 13 kDa ribonucleoprotein-associated protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	3G	121	Total	C	N	O	S	0	0
			916	583	158	171	4		
24	3H	121	Total	C	N	O	S	0	0
			916	583	158	171	4		

- Molecule 25 is a protein called U3 small nucleolar RNA-associated protein 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	A4	662	Total	C	N	O	S	0	0
			5226	3309	910	986	21		

- Molecule 26 is a protein called U3 small nucleolar RNA-associated protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	A5	514	Total	C	N	O	S	0	0
			3976	2520	688	755	13		

- Molecule 27 is a protein called U3 small nucleolar RNA-associated protein 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	A8	548	Total	C	N	O	S	0	0
			3307	2054	608	642	3		

- Molecule 28 is a protein called U3 small nucleolar RNA-associated protein 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	A9	128	Total	C	N	O	S	0	0
			939	594	173	170	2		

- Molecule 29 is a protein called U3 small nucleolar RNA-associated protein 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	AE	1534	Total	C	N	O	S	0	0
			9955	6242	1771	1923	19		

- Molecule 30 is a protein called U3 small nucleolar RNA-associated protein 15.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	AF	493	Total	C	N	O	S	0	0
			3911	2462	702	735	12		

- Molecule 31 is a protein called NET1-associated nuclear protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	AG	826	Total	C	N	O	S	0	0
			6570	4181	1111	1259	19		

- Molecule 32 is a protein called Periodic tryptophan protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	B1	834	Total	C	N	O	S	0	0
			6635	4223	1140	1253	19		

- Molecule 33 is a protein called U3 small nucleolar RNA-associated protein 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	B2	851	Total	C	N	O	S	0	0
			6723	4294	1133	1269	27		

- Molecule 34 is a protein called U3 small nucleolar RNA-associated protein 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	B3	757	Total	C	N	O	S	0	0
			5919	3769	993	1130	27		

- Molecule 35 is a protein called U3 small nucleolar RNA-associated protein 18.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	B8	477	Total	C	N	O	S	0	0
			3764	2387	662	705	10		

- Molecule 36 is a protein called U3 small nucleolar RNA-associated protein 21.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	BE	865	Total	C	N	O	S	0	0
			6810	4322	1175	1292	21		

- Molecule 37 is a protein called U3 small nucleolar RNA-associated protein 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	B6	374	Total	C	N	O	S	0	0
			2800	1782	501	505	12		

- Molecule 38 is a protein called Bud site selection protein 21.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	5B	60	Total	C	N	O		0	0
			495	310	101	84			

- Molecule 39 is a protein called U3 small nucleolar RNA-associated protein 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	5C	535	Total	C	N	O	S	0	0
			4237	2656	762	807	12		

- Molecule 40 is a protein called U3 small nucleolar RNA-associated protein 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	5D	167	Total	C	N	O	S	0	0
			1396	862	266	263	5		

- Molecule 41 is a protein called U3 small nucleolar RNA-associated protein MPP10.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	5E	204	Total	C	N	O	S	0	0
			1647	1021	294	328	4		

- Molecule 42 is a protein called U3 small nucleolar ribonucleoprotein protein IMP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	5F	182	Total	C	N	O	S	0	0
			1530	967	287	269	7		

- Molecule 43 is a protein called U3 small nucleolar ribonucleoprotein protein IMP4.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	5G	282	Total	C	N	O	S	0	0
			2296	1441	430	418	7		

- Molecule 44 is a protein called Something about silencing protein 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	5H	74	Total	C	N	O	S	0	0
			596	373	122	101			

- Molecule 45 is a protein called Protein SOF1.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	5I	461	Total	C	N	O	S	0	0
			3765	2354	686	709	16		

- Molecule 46 is a protein called rRNA-processing protein FCF2.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	5J	151	Total	C	N	O	S	0	0
			1280	807	240	228	5		

- Molecule 47 is a protein called rRNA-processing protein FCF1.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	5K	175	Total	C	N	O	S	0	0
			1403	896	256	241	10		

- Molecule 48 is a protein called Ribosome biogenesis protein ENP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	RA	338	Total	C	N	O	S	0	0
			2709	1713	463	524	9		

- Molecule 49 is a protein called U3 small nucleolar ribonucleoprotein protein LCP5.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	RB	134	Total	C	N	O	S	0	0
			1108	664	227	214	3		

- Molecule 50 is a protein called KRR1 small subunit processome component.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	RC	278	Total	C	N	O	S	0	0
			2207	1408	391	395	13		

- Molecule 51 is a protein called rRNA biogenesis protein RRP5.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	RD	316	Total	C	N	O	S	0	0
			2412	1541	414	452	5		

- Molecule 52 is a protein called U3 small nucleolar RNA-associated protein 22.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	RE	1079	Total	C	N	O	S	0	0
			8716	5666	1437	1589	24		

- Molecule 53 is a protein called Ribosomal RNA-processing protein 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	RF	241	Total	C	N	O	S	0	0
			1963	1253	335	367	8		

- Molecule 54 is a protein called Ribosomal RNA small subunit methyltransferase NEP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	RG	216	Total	C	N	O	S	0	0
			1701	1079	296	315	11		
54	RH	230	Total	C	N	O	S	0	0
			1799	1142	313	333	11		

- Molecule 55 is a protein called Ribosome biogenesis protein BMS1.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	RJ	796	Total	C	N	O	S	0	0
			6379	4086	1136	1128	29		

- Molecule 56 is a protein called RNA 3'-terminal phosphate cyclase-like protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	RK	360	Total	C	N	O	S	0	0
			2781	1781	473	516	11		

- Molecule 57 is a protein called RNA cytidine acetyltransferase.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	RL	805	Total	C	N	O	S	0	0
			4539	2760	885	887	7		
57	RM	766	Total	C	N	O		0	0
			3779	2247	766	766			

- Molecule 58 is a protein called Nucleolar complex protein 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	RN	607	Total	C	N	O	S	0	0
			4529	2861	820	837	11		

- Molecule 59 is a protein called Nucleolar complex protein 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	RO	525	Total	C	N	O	S	0	0
			3766	2412	646	696	12		

- Molecule 60 is a protein called U3 small nucleolar RNA-associated protein 20.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	RP	2108	Total	C	N	O	S	0	0
			12171	7483	2291	2381	16		

- Molecule 61 is a protein called U3 small nucleolar RNA-associated protein 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	RQ	226	Total	C	N	O	S	0	0
			1651	1023	313	313	2		

- Molecule 62 is a protein called Essential nuclear protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	RS	251	Total	C	N	O	S	0	0
			2051	1340	349	359	3		

- Molecule 63 is a protein called Pno1.

Mol	Chain	Residues	Atoms					AltConf	Trace
63	RT	171	Total	C	N	O	S	0	0
			1357	864	249	240	4		

- Molecule 64 is a protein called Protein FAF1.

Mol	Chain	Residues	Atoms					AltConf	Trace
64	RV	190	Total	C	N	O	S	0	0
			1448	891	290	264	3		

- Molecule 65 is a protein called Protein BFR2.

Mol	Chain	Residues	Atoms				AltConf	Trace
65	RY	37	Total	C	N	O	0	0
			299	191	48	60		

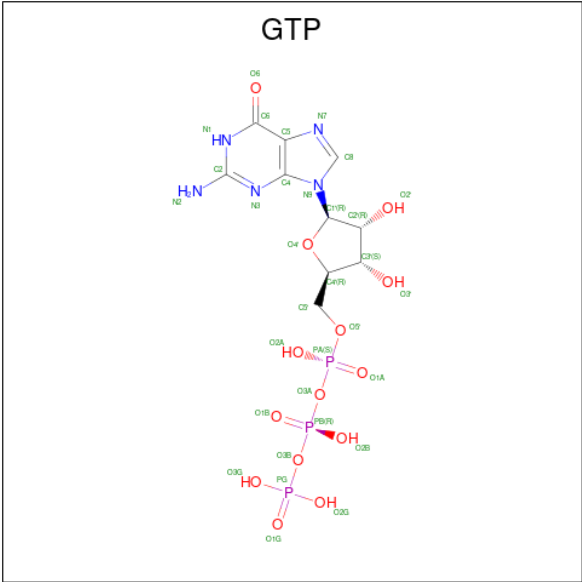
- Molecule 66 is a protein called Unassigned helices.

Mol	Chain	Residues	Atoms				AltConf	Trace
66	X1	75	Total	C	N	O	0	0
			375	225	75	75		

- Molecule 67 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
67	Sc	1	Total	Zn	0
			1	1	
67	5K	1	Total	Zn	0
			1	1	

- Molecule 68 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					AltConf
68	RJ	1	Total	C	N	O	P	0
			32	10	5	14	3	

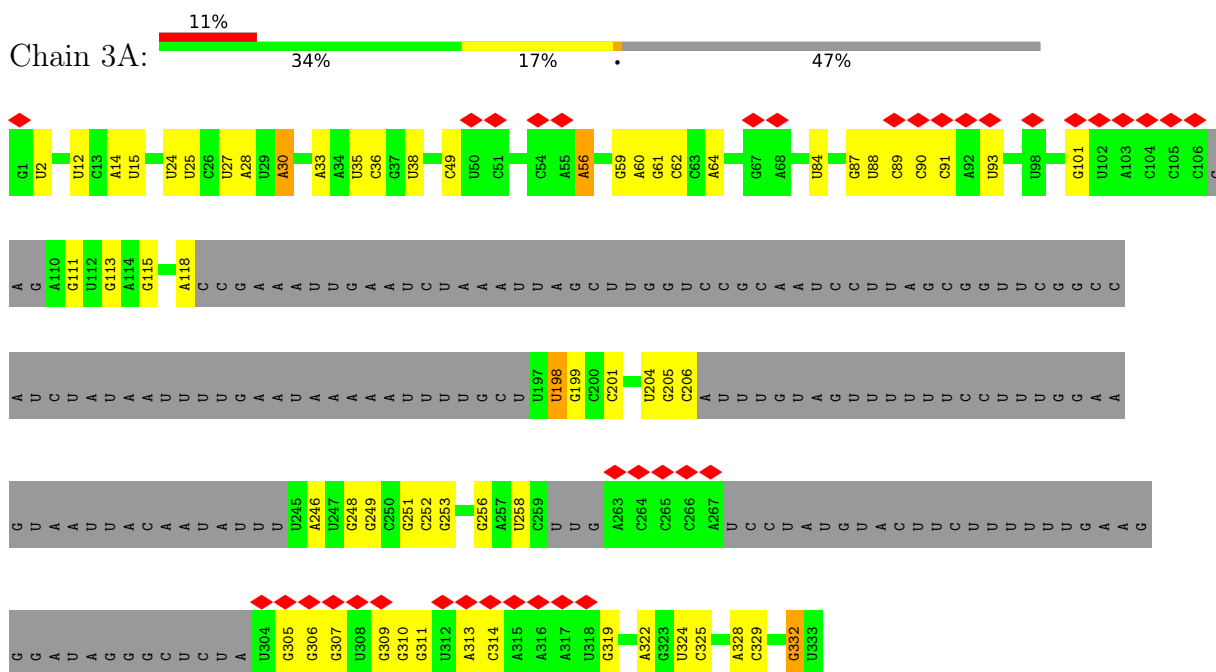
- Molecule 69 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
69	RJ	1	Total	Mg	0
			1	1	

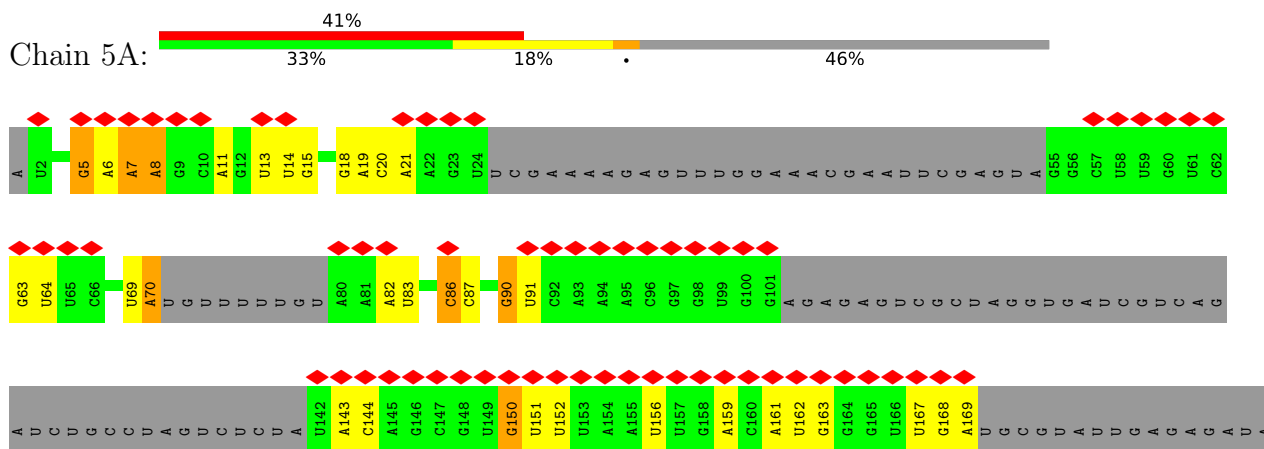
3 Residue-property plots [i](#)

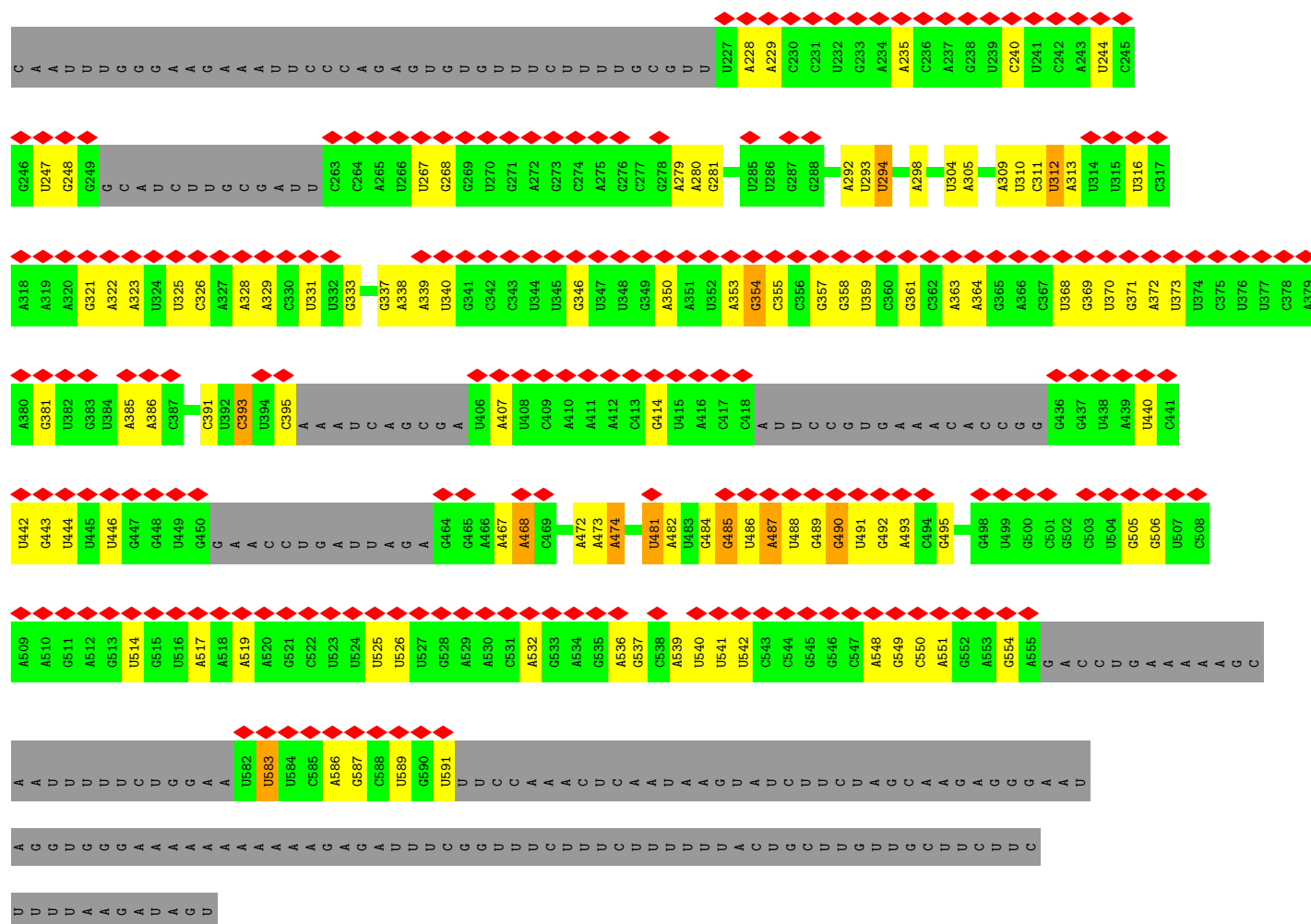
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: U3 snoRNA

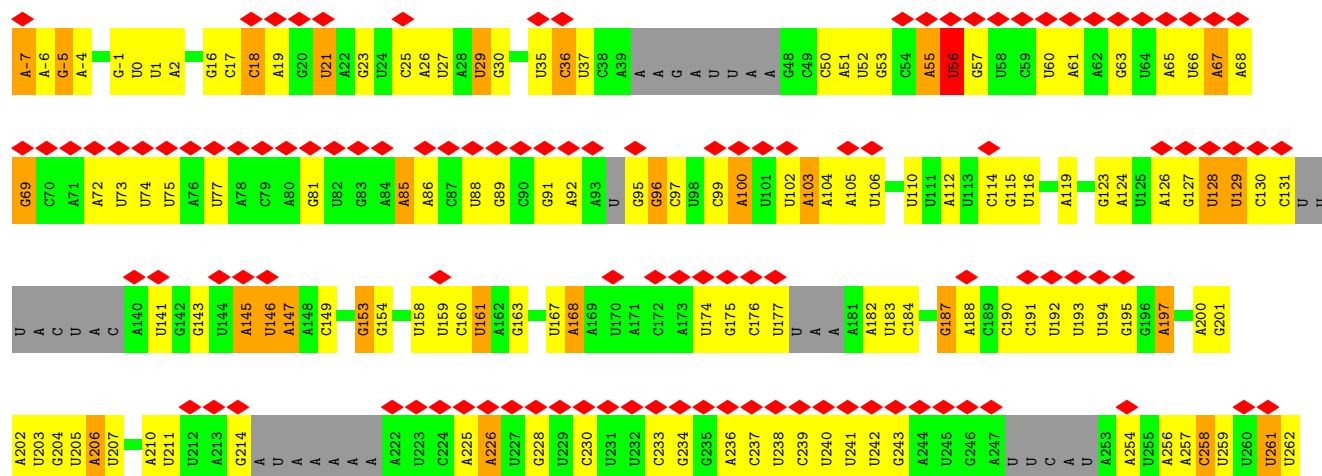


• Molecule 2: 5' ETS

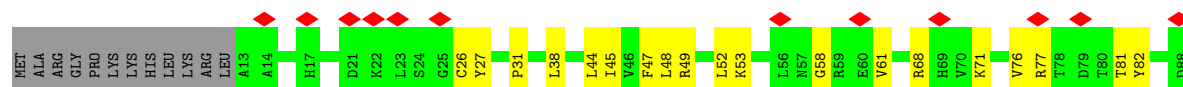


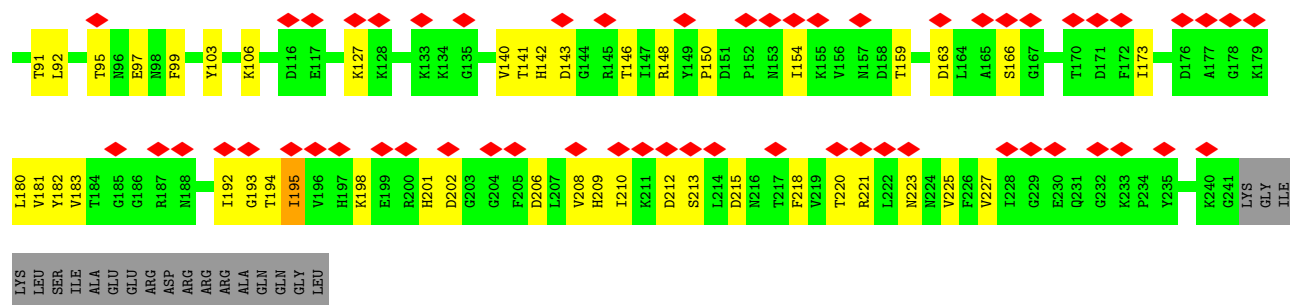


• Molecule 3: 18S pre-rRNA

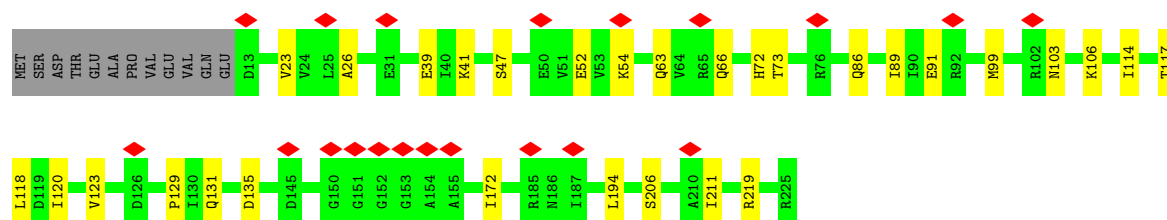
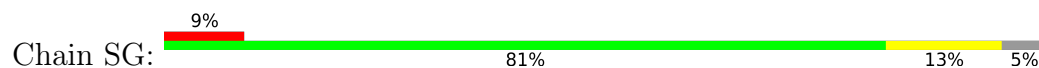




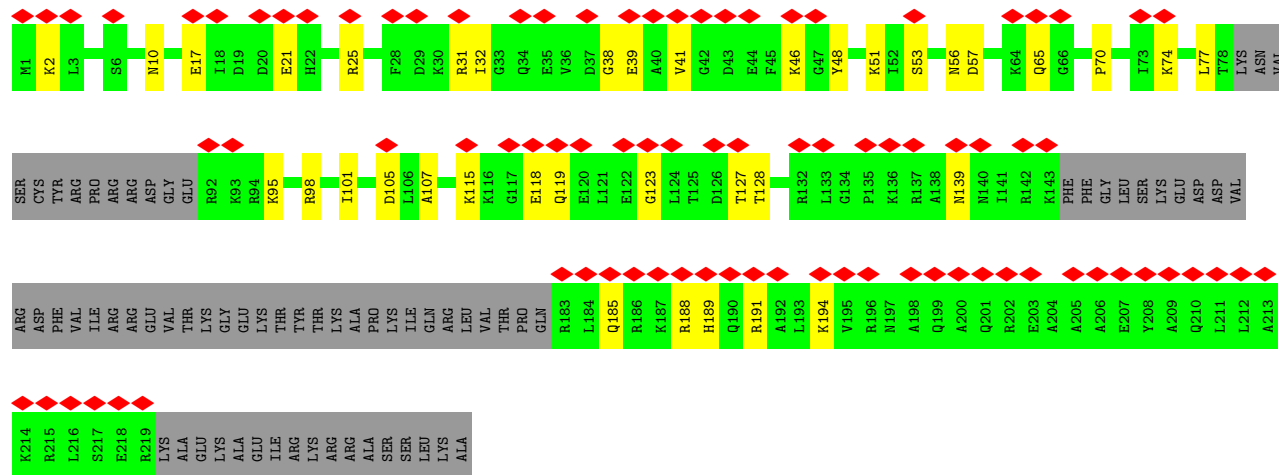




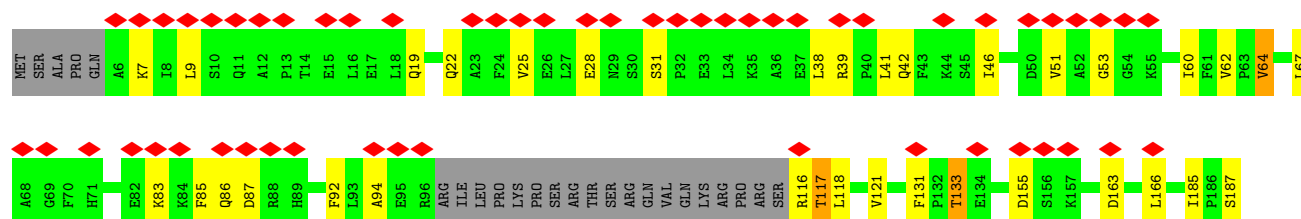
• Molecule 6: 40S ribosomal protein S5



• Molecule 7: 40S ribosomal protein S6-A

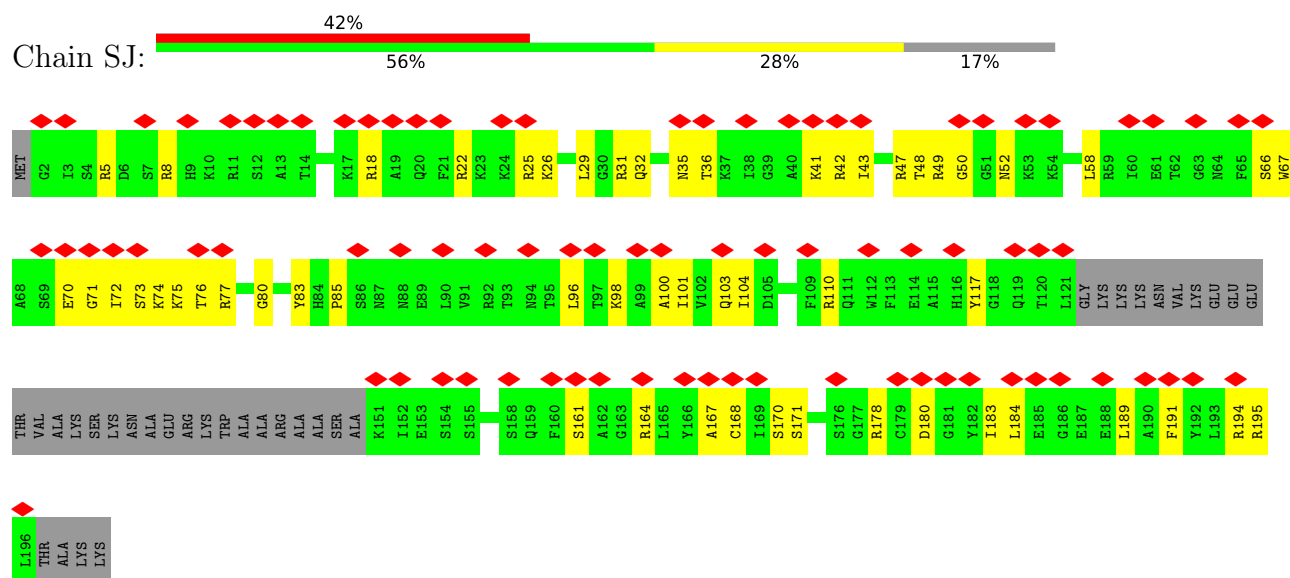


• Molecule 8: 40S ribosomal protein S7-A

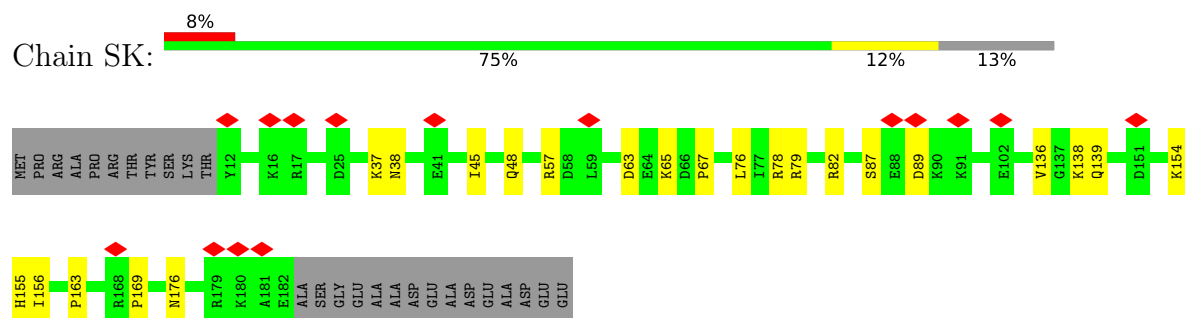




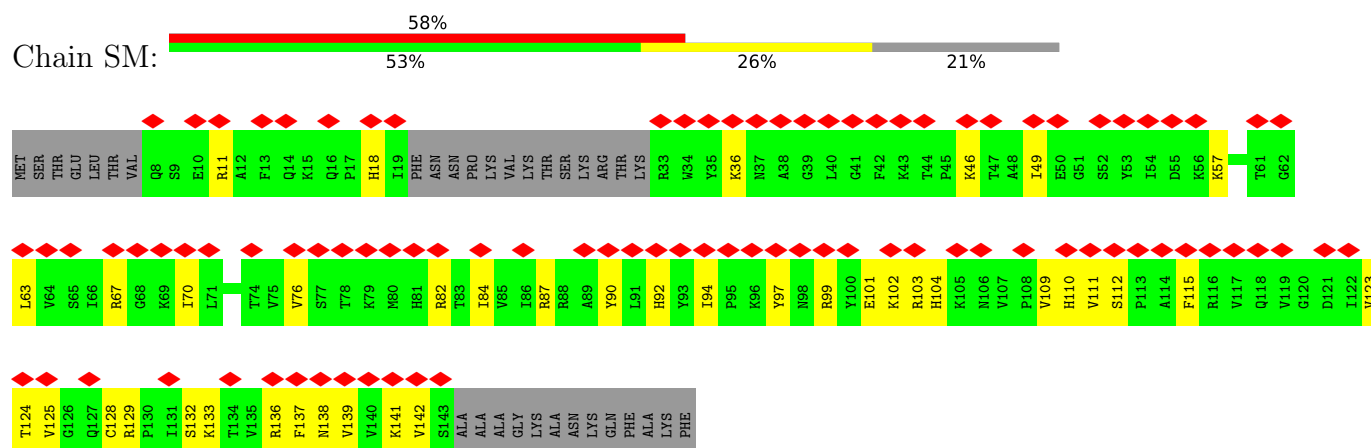
• Molecule 9: 40S ribosomal protein S8-A



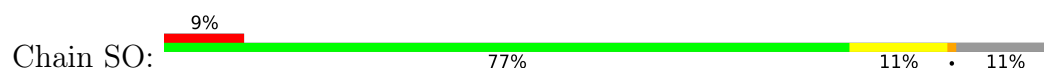
• Molecule 10: 40S ribosomal protein S9-A



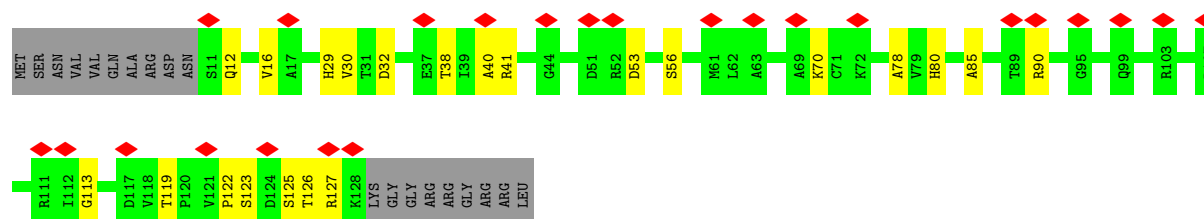
• Molecule 11: 40S ribosomal protein S11-A



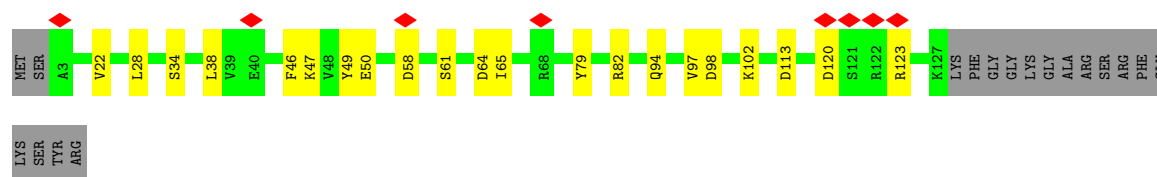
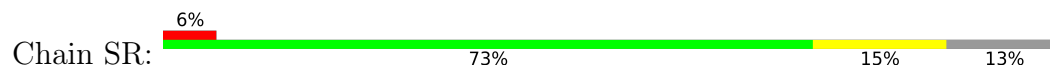
• Molecule 12: 40S ribosomal protein S13



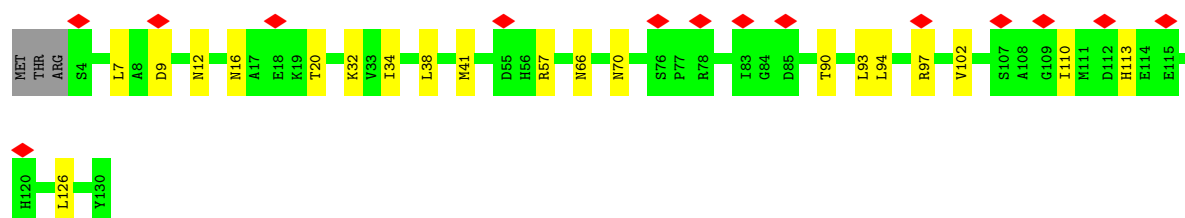
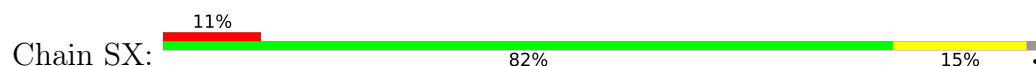
• Molecule 13: 40S ribosomal protein S14-A



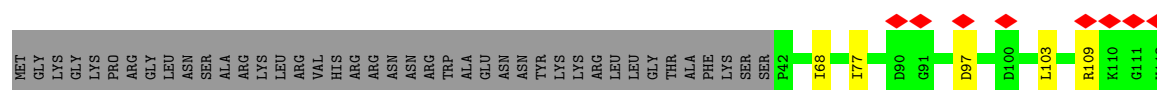
• Molecule 14: 40S ribosomal protein S16-A



• Molecule 15: 40S ribosomal protein S22-B

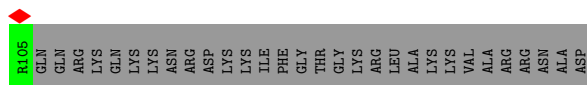


• Molecule 16: 40S ribosomal protein S23-A

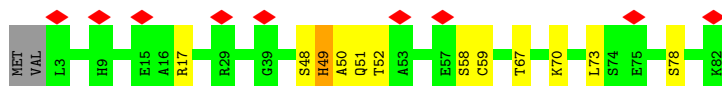
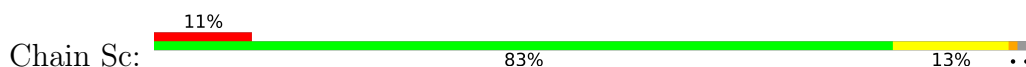




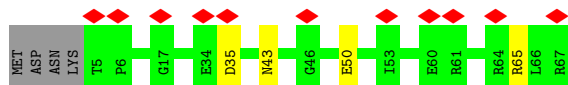
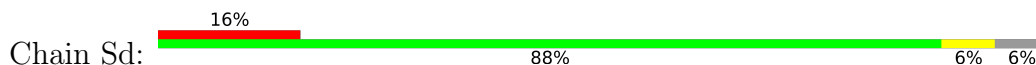
- Molecule 17: 40S ribosomal protein S24-A



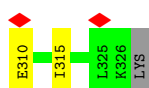
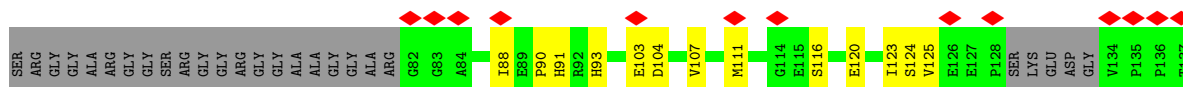
- Molecule 18: 40S ribosomal protein S27-A



- Molecule 19: 40S ribosomal protein S28-A

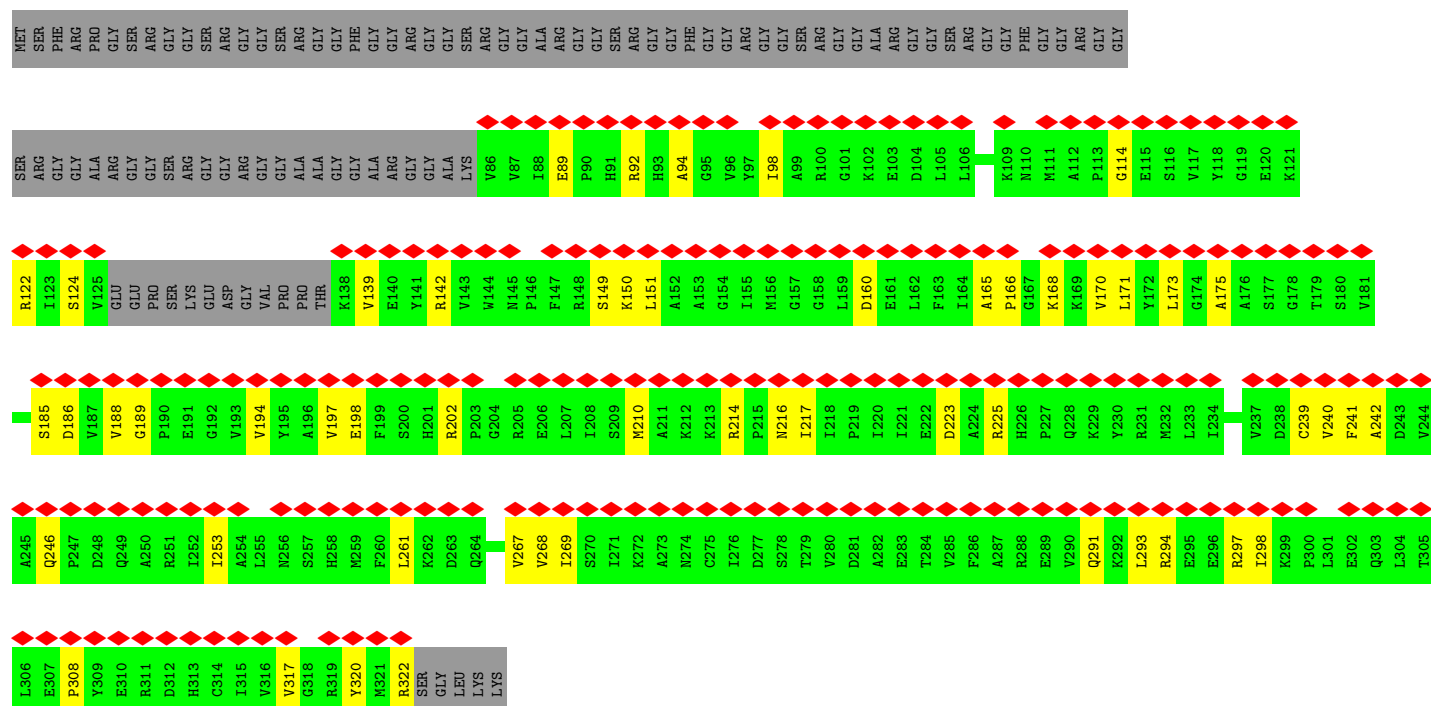


- Molecule 20: rRNA 2'-O-methyltransferase fibrillar

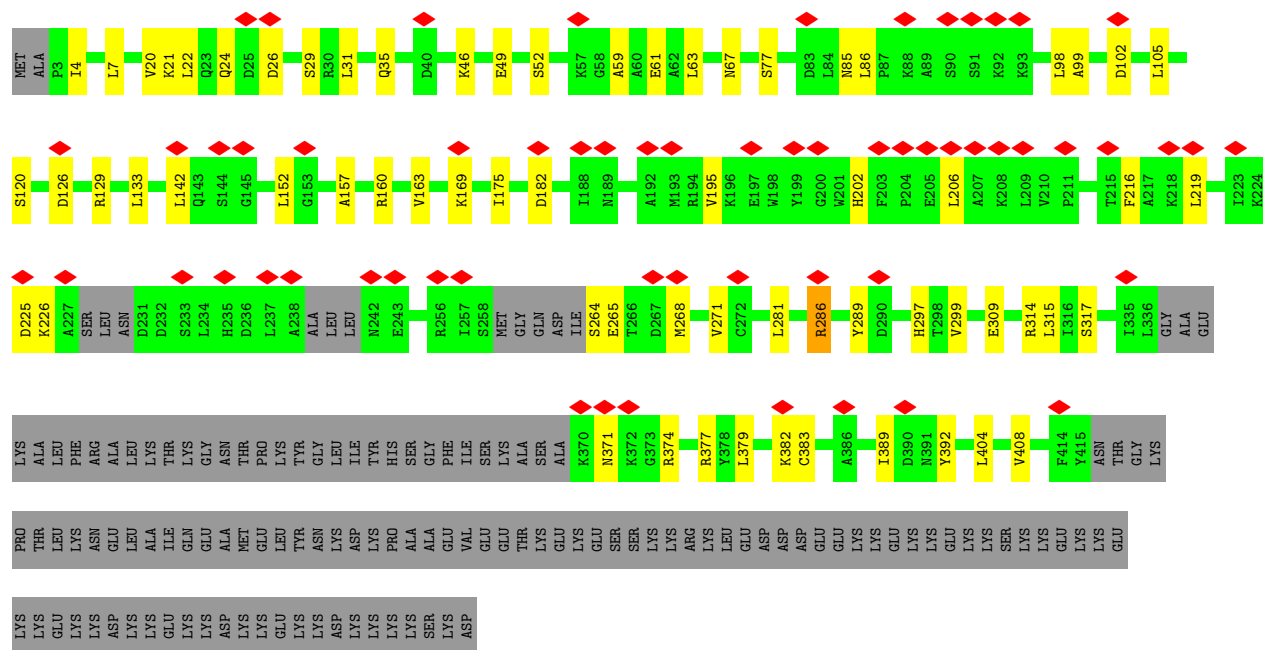


- Molecule 20: rRNA 2'-O-methyltransferase fibrillar

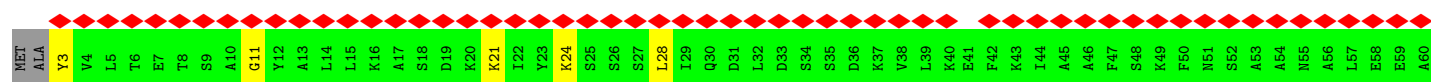


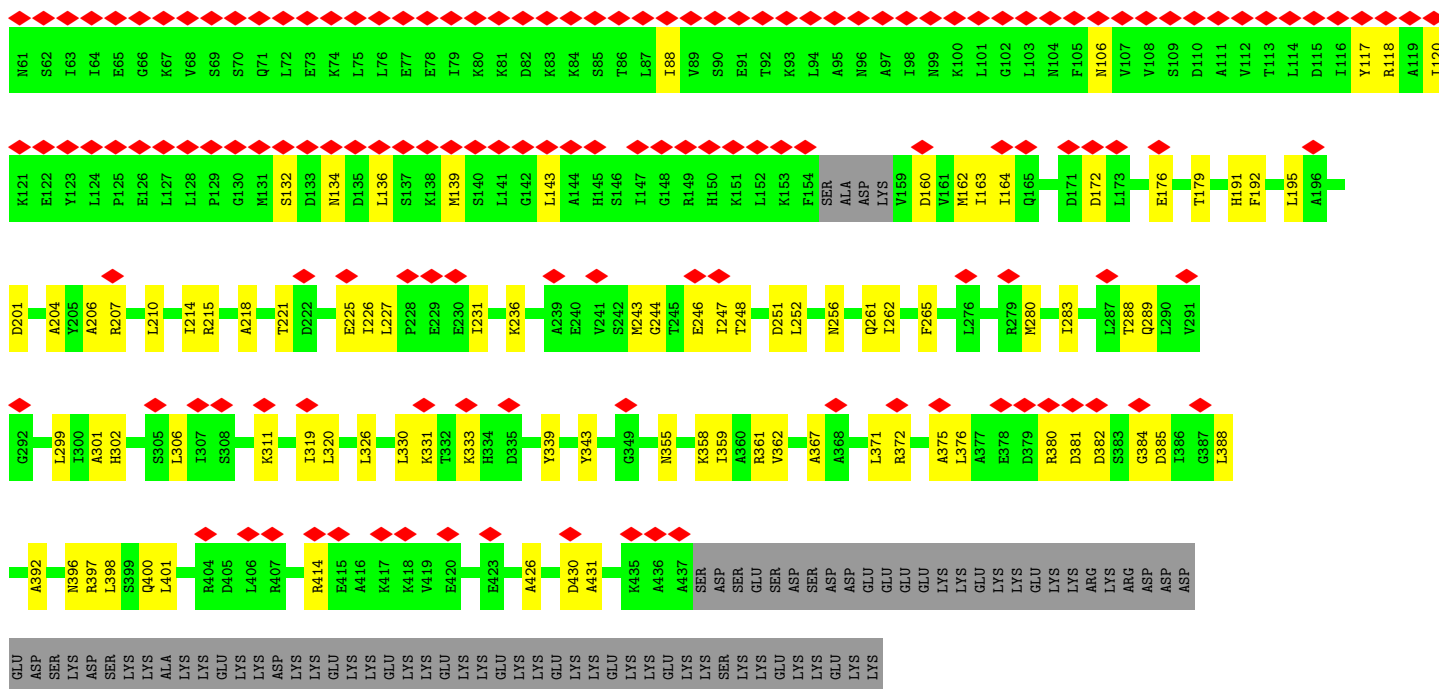


• Molecule 21: Nucleolar protein 56

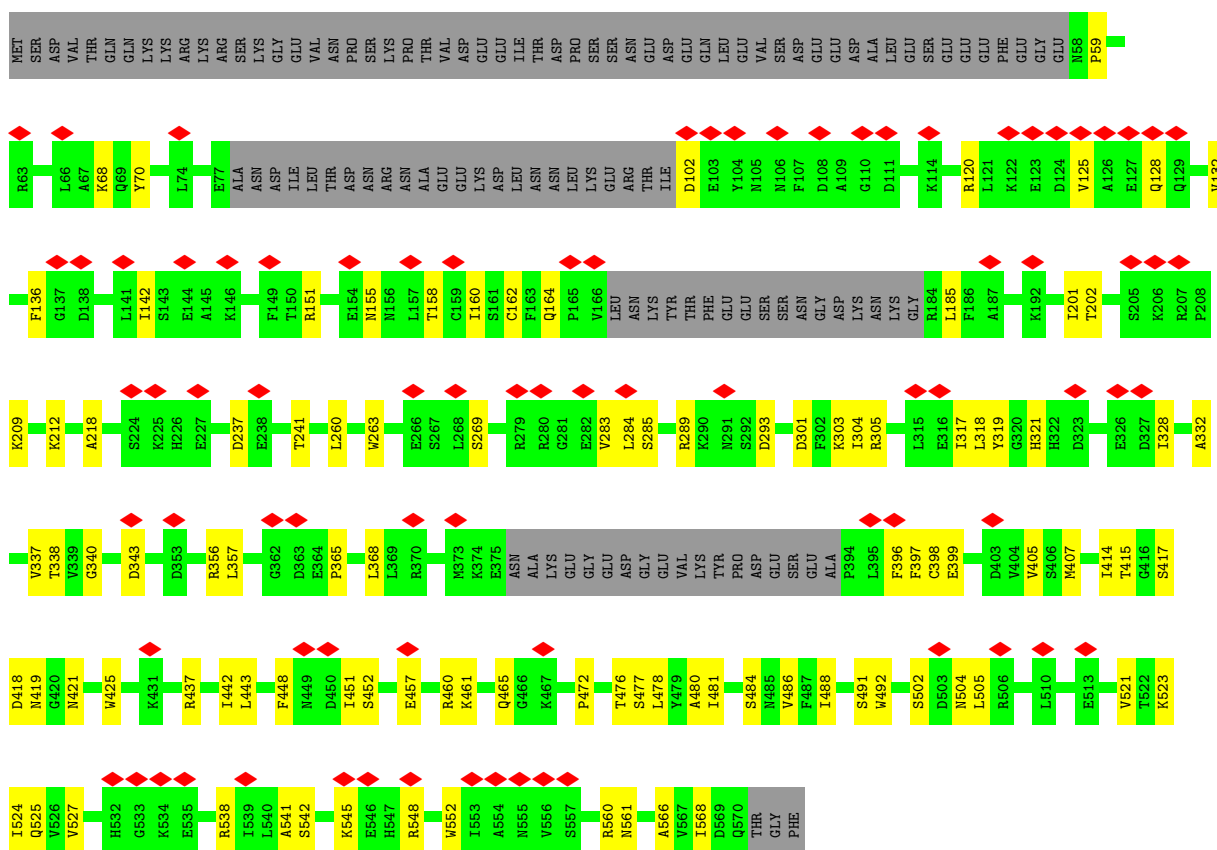


• Molecule 22: Nucleolar protein 58

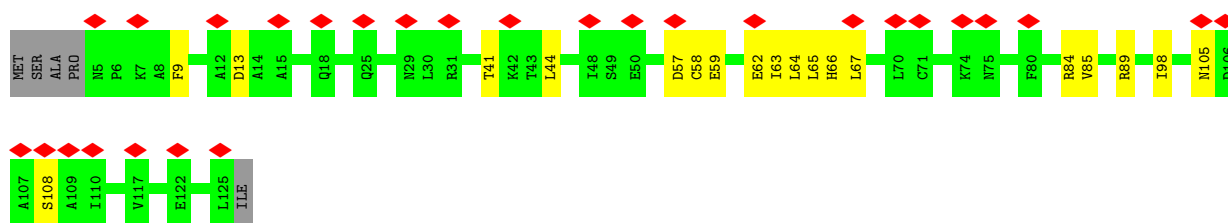
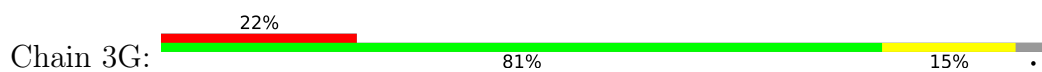




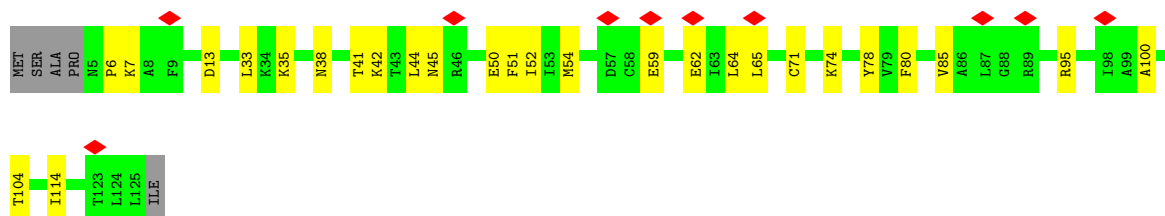
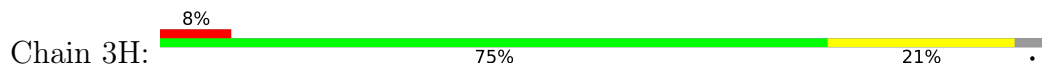
• Molecule 23: Ribosomal RNA-processing protein 9



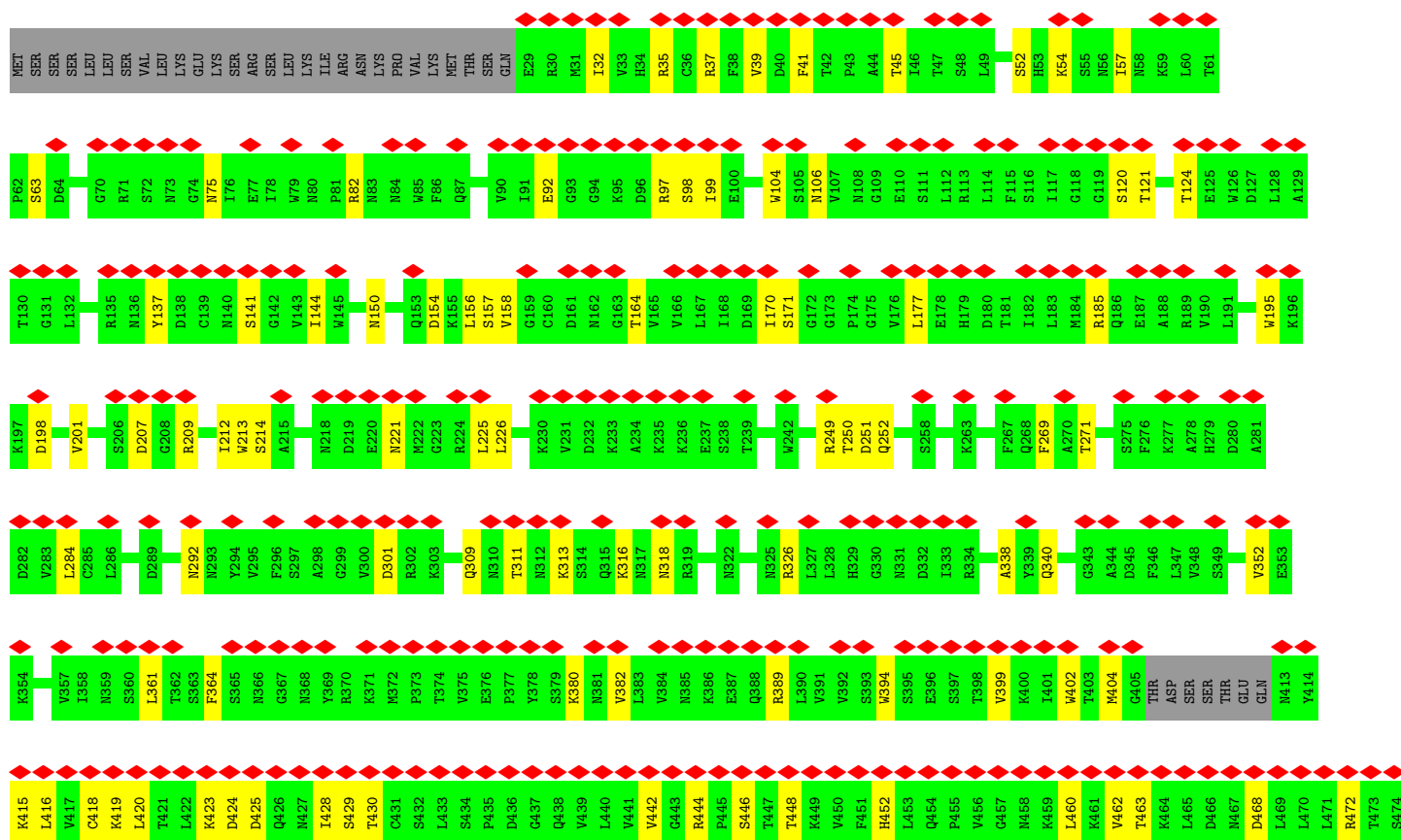
• Molecule 24: 13 kDa ribonucleoprotein-associated protein

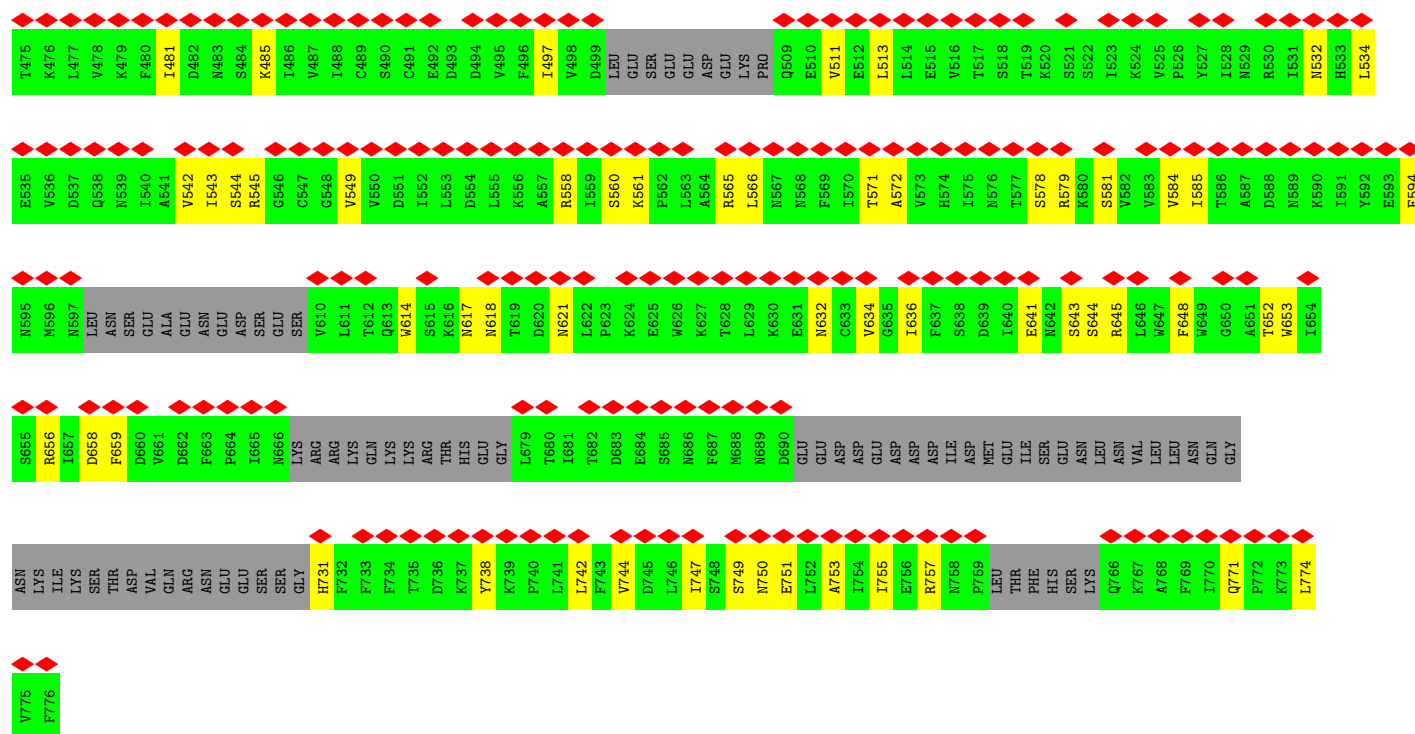


- Molecule 24: 13 kDa ribonucleoprotein-associated protein

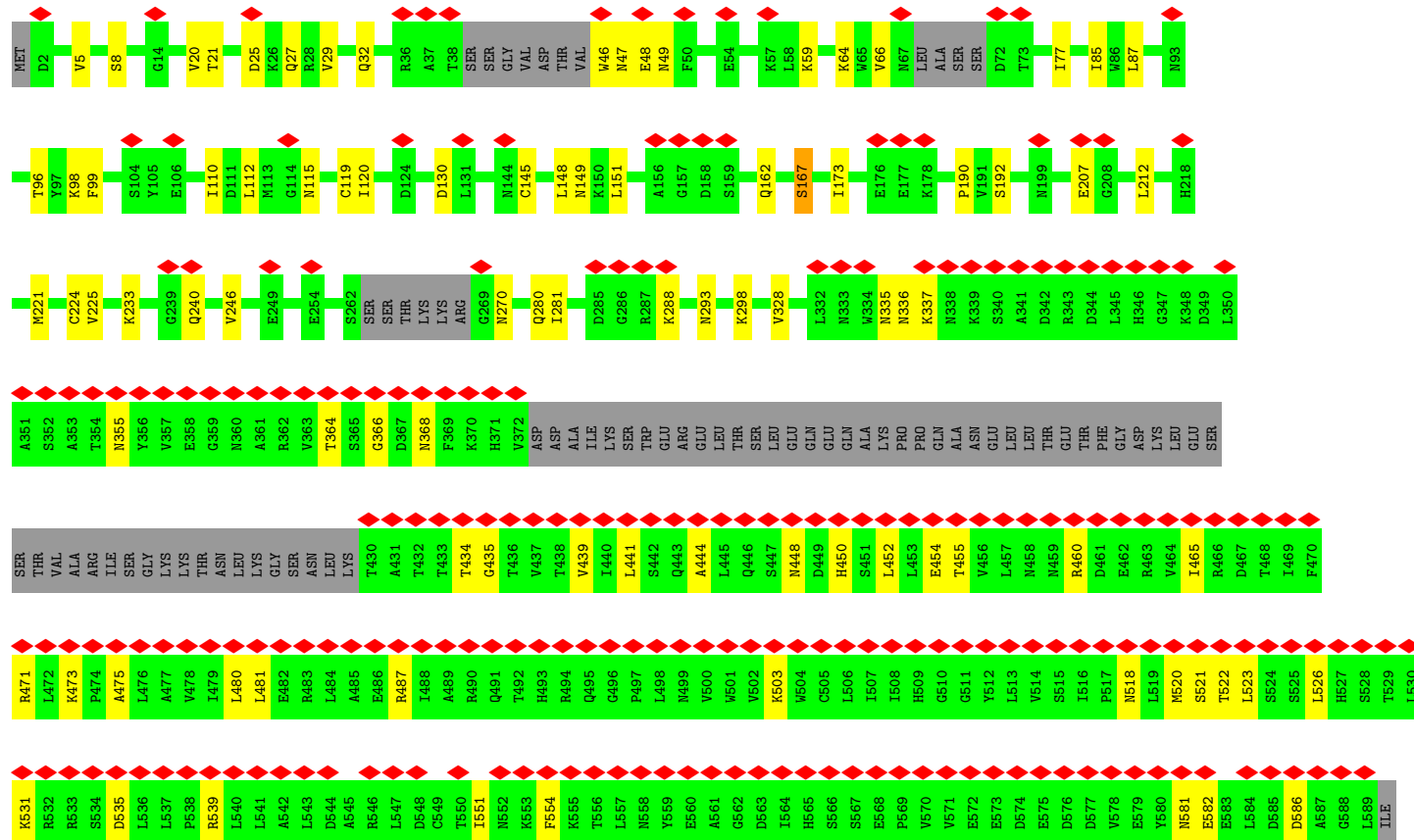
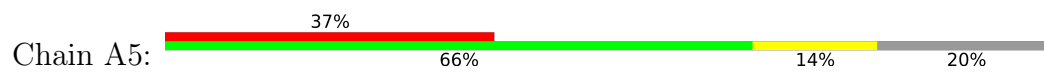


- Molecule 25: U3 small nucleolar RNA-associated protein 4





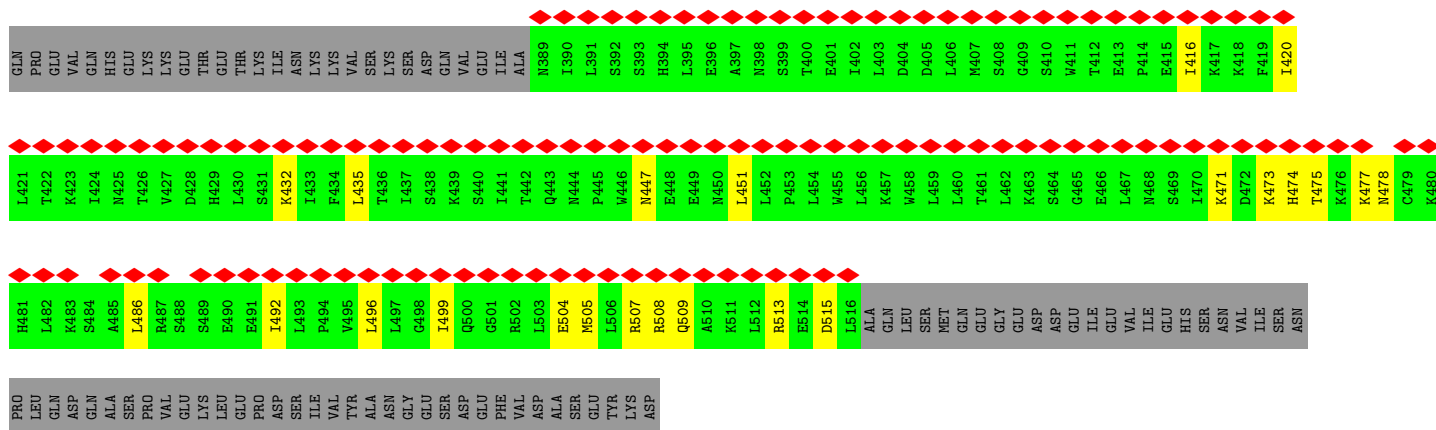
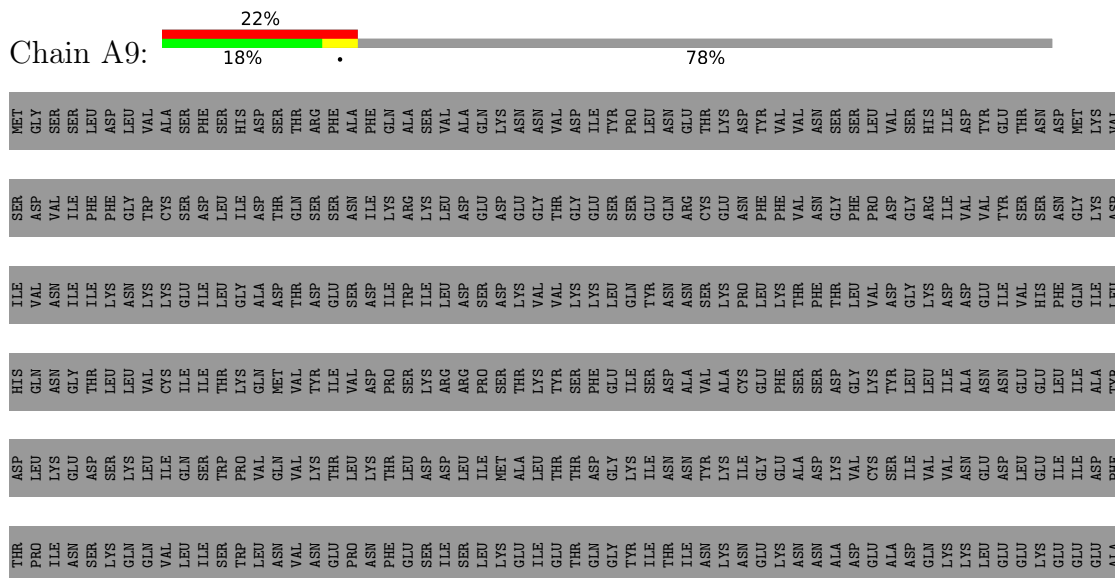
• Molecule 26: U3 small nucleolar RNA-associated protein 5



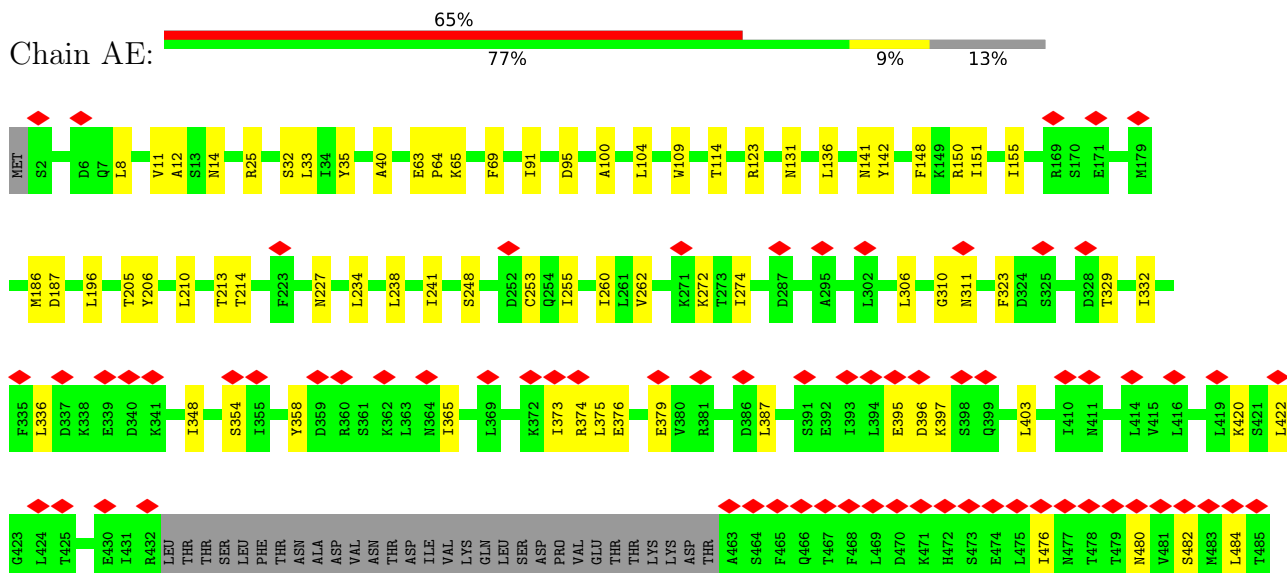
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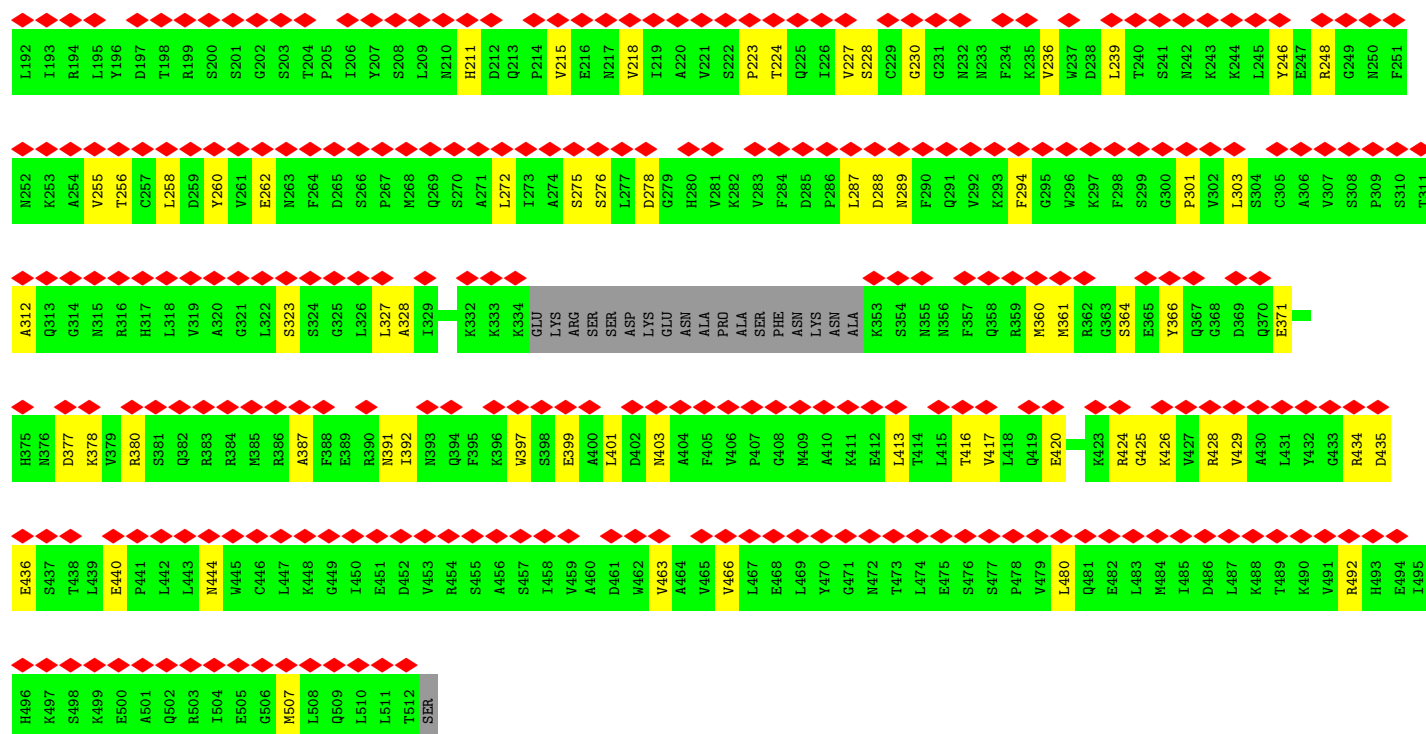
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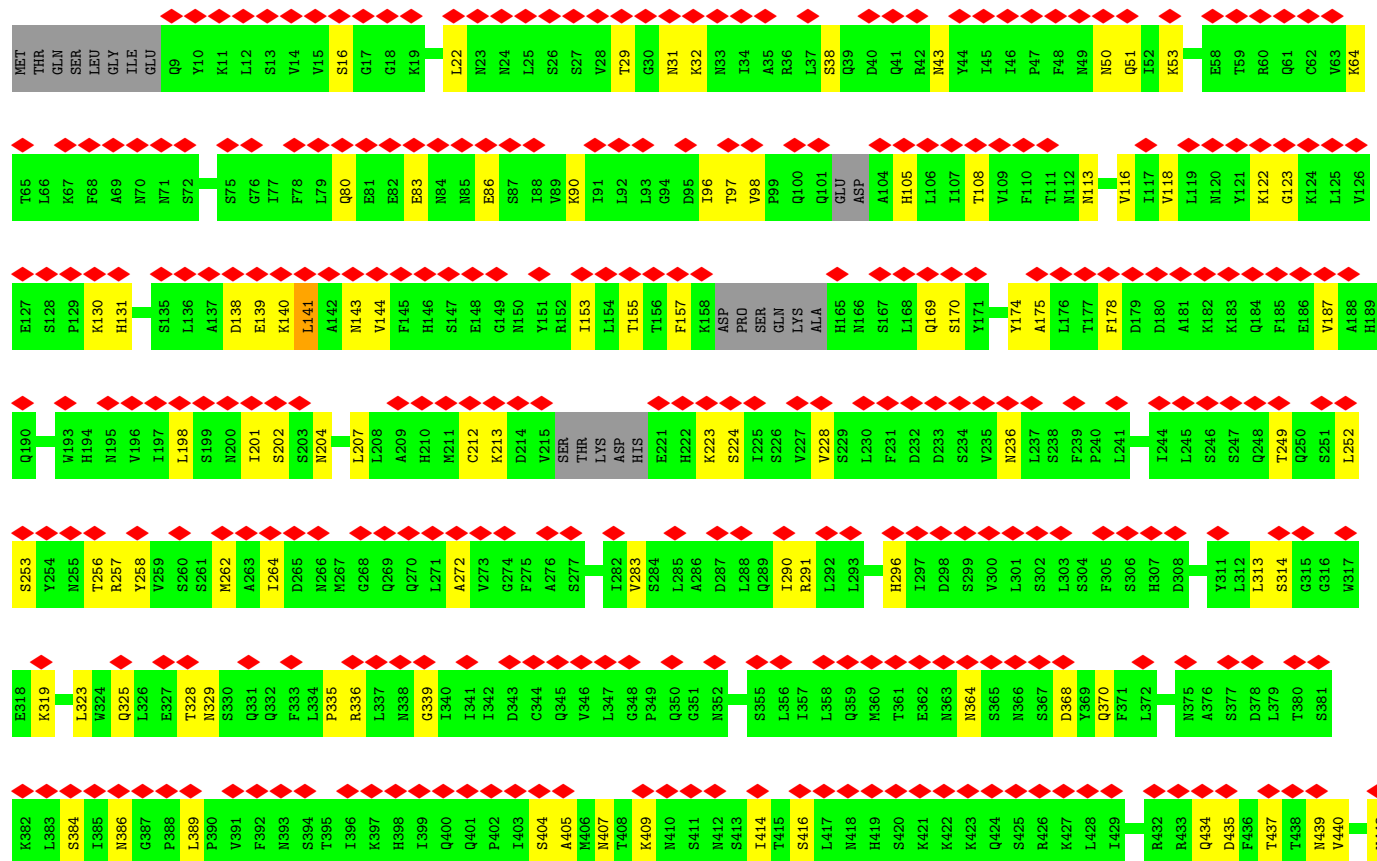
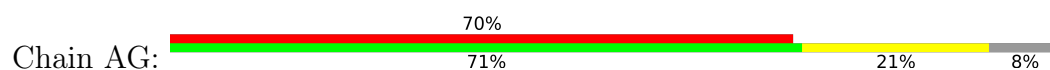
Chain AE:



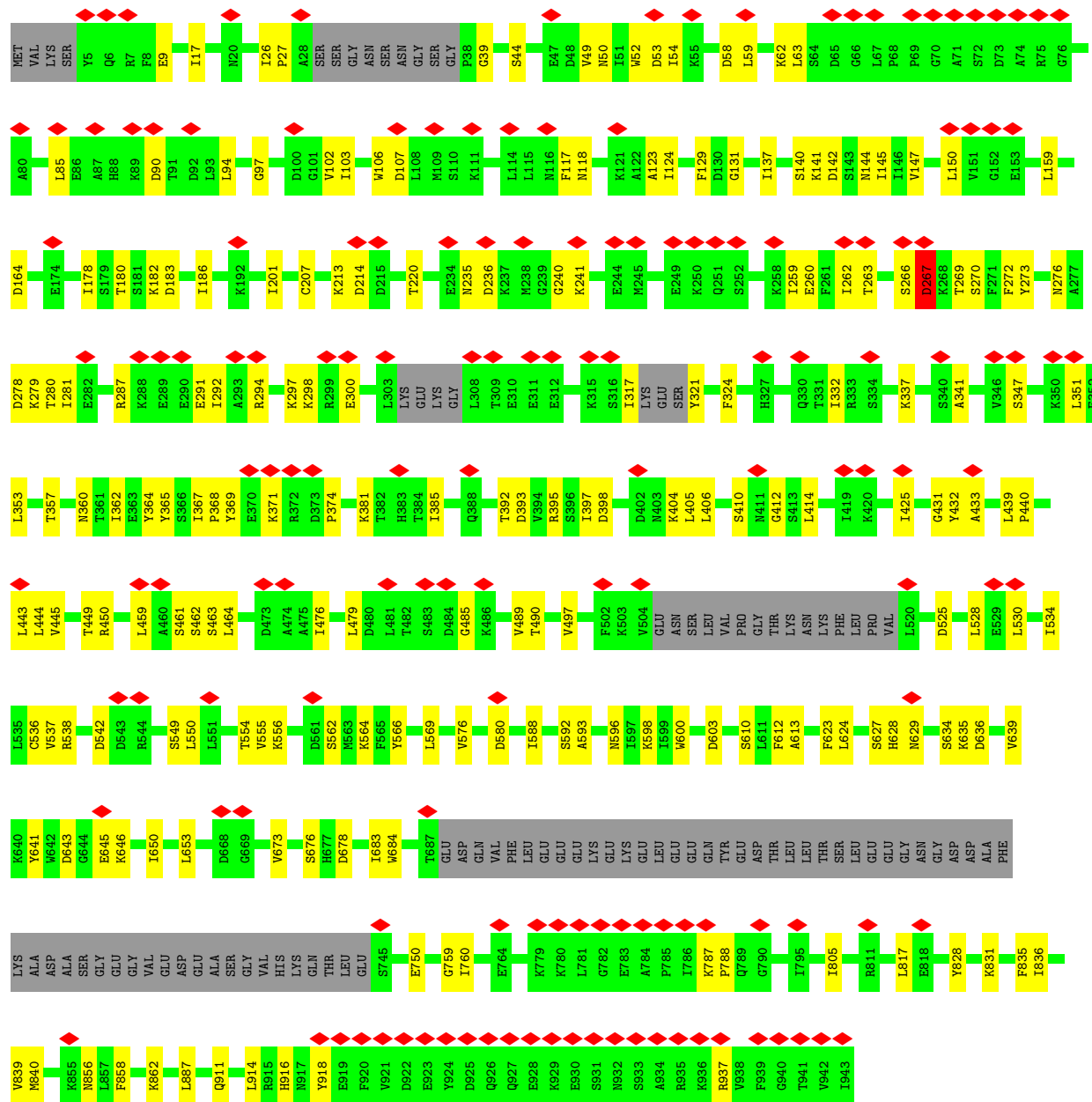
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GLU	THR	ASP	GLN	D1150	Y1151	Y1152	D1153	V1154	R1155	K1156	M1157	L1158	R1159	L1160	K1161	V1162	Y1163	S1164	V1165	L1166	LEU	ASP	GLU	THR	SER	D1172	K1173	K1174	L1175	I1176	R1177	M1178	I1179	R1180	GLU	GLU	PHE	GLY	THR	LEU	LEU	GLU	GLY	VAL	LEU	PHE	PHE	ILE	ASN	SER	V1197	E1198	L1199	T1200	F1201	S1202	C1203	I1204	T1205	
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V1026	K1027	L1028	F1029	S1030	T1031	L1032	ILE	LYS	THR	ASP	PRO	V1039	K1040	A1041	L1042	G1043	S1044	F1045	L1046	F1047	L1048	I1049	ALA	GLN	TYR	SER	SER	ALA	L1057	V1058	N1059	F1060	K1061	I1062	G1063	E1064	A1065	R1066	L1067	L1068	L1069	E1070	F1071	ILE	LYS	ALA	LEU	VAL	ASP	LEU	H1080	V1081	M1082	E1083	L1084	L1085				
F966	M967	G968	ALA	HIS	SER	ILE	ARG	GLN	ASP	GLU	THR	THR	LYS	VAL	VAL	GLU	ARG	THR	VAL	VAL	P992	A993	L994	I995	K996	N997	S998	K999	G1000	N1001	E1002	K1003	E1004	E1005	M1006	GLU	PHE	LEU	LEU	LEU	SER	PHE	THR	THR	ALA	LEU	LEU	GLN	H1019	V1020	P1021	R1022	H1023	R1024	N1025					
ILE	THR	Y908	L909	K910	H911	H912	G913	C914	T915	E916	L917	T918	N919	V920	R921	ALA	ASP	ILE	VAL	SER	ALA	ILE	ARG	ASN	SER	A933	S934	P935	Q936	Y937	Q938	N939	K940	L941	L942	L943	V944	I945	G946	S947	L948	ALA	THR	THR	LEU	SER	SER	E954	V955	L956	L957	H958	S959	V960	N961	P962	I963	F964	T965	
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P726	K727	E728	K729	Q730	S731	F732	M733	I734	D735	F736	V737	L738	S739	A740	L741	N742	S743	D744	V745	E746	Q747	L748	A749	N750	I751	A752	A753	E754	Y755	L756	I757	S758	F759	F760	A761	S762	L763	N764	N765	A766	Q767	K768	L769	K770	I771	V772	Q773	N774	I775	V776	D777	S778	S779	S780	N781	V782	E783	S784	S785	
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• Molecule 31: NET1-associated nuclear protein 1

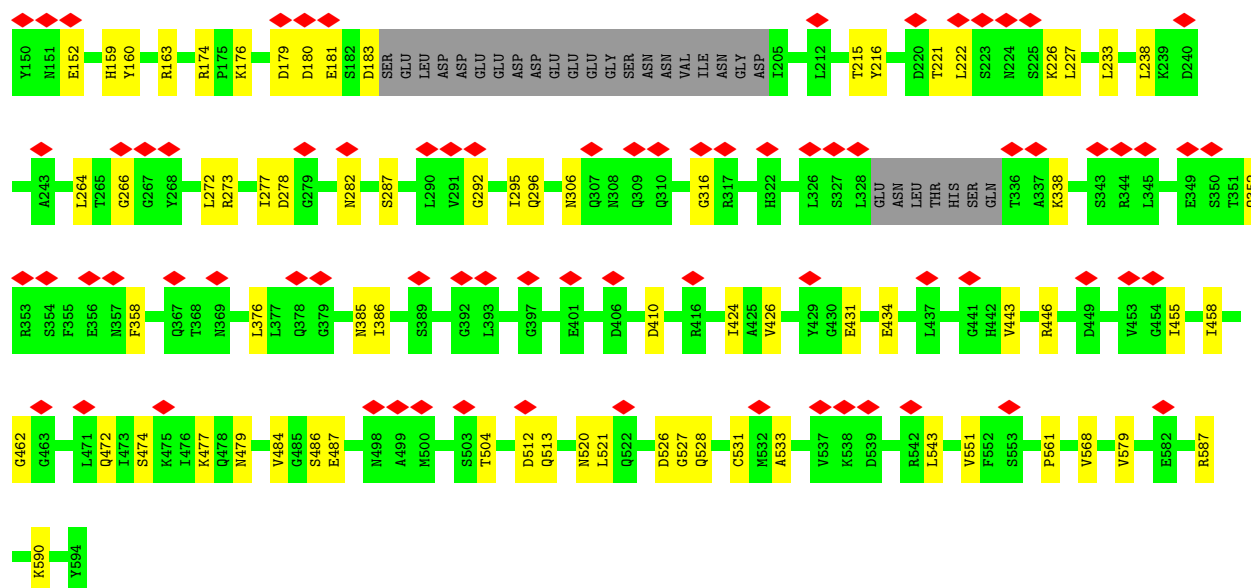


- Molecule 33: U3 small nucleolar RNA-associated protein 12

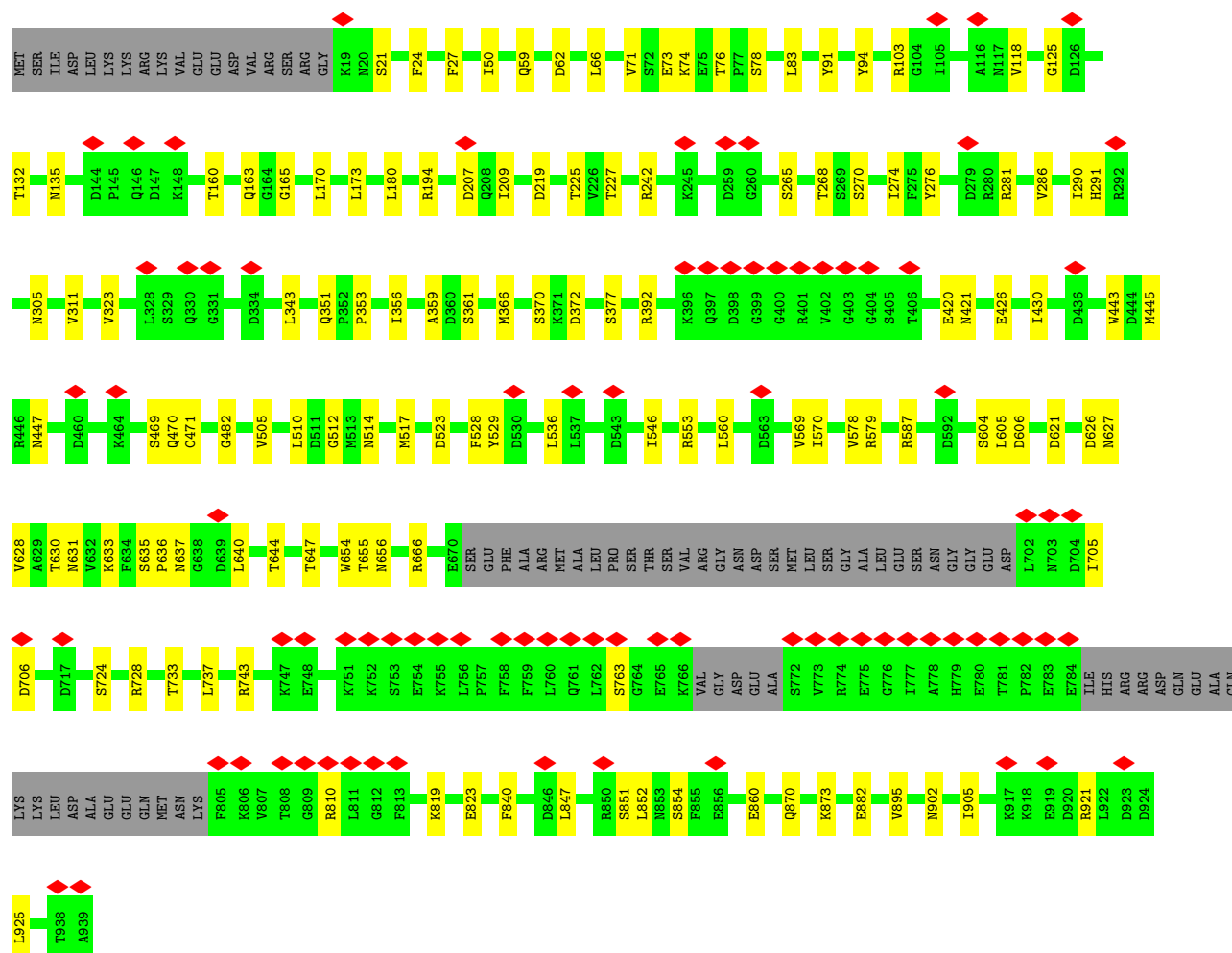
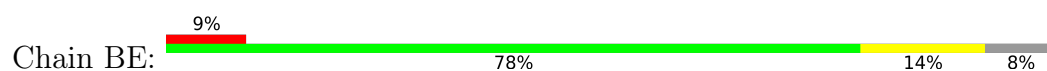


- Molecule 34: U3 small nucleolar RNA-associated protein 13



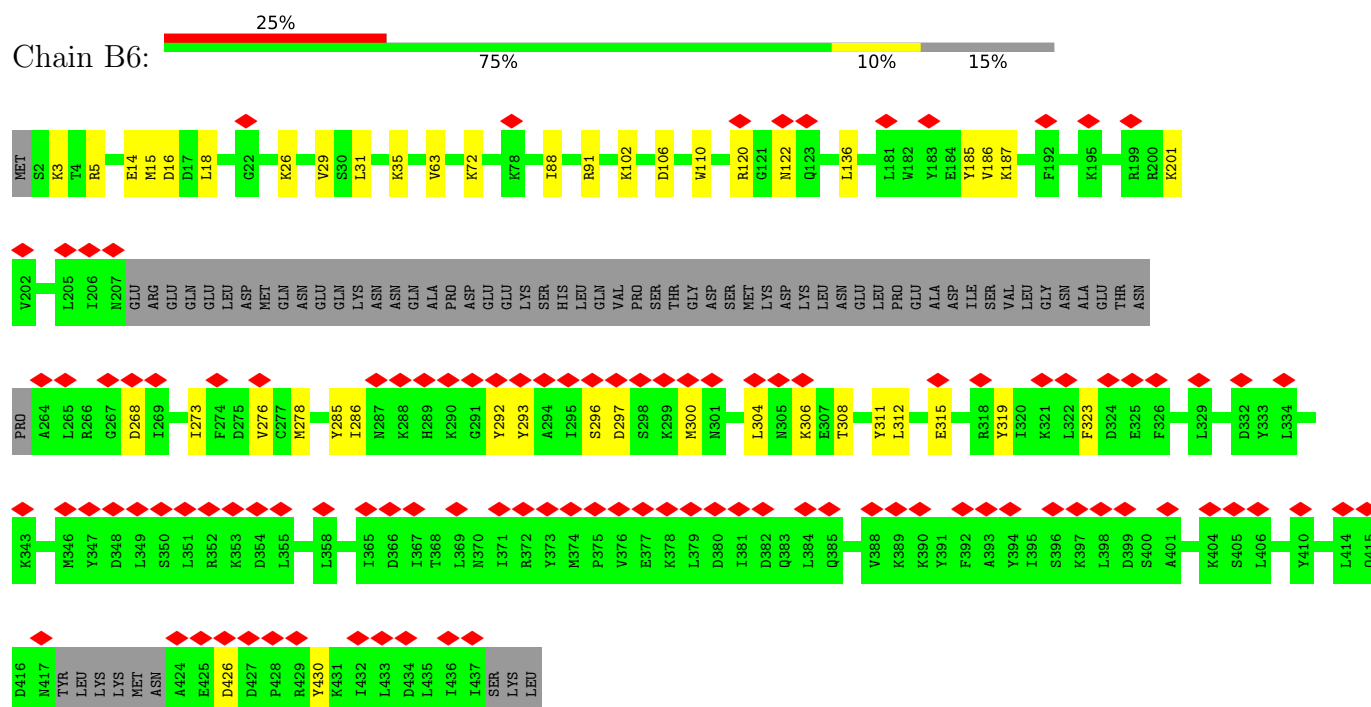


• Molecule 36: U3 small nucleolar RNA-associated protein 21



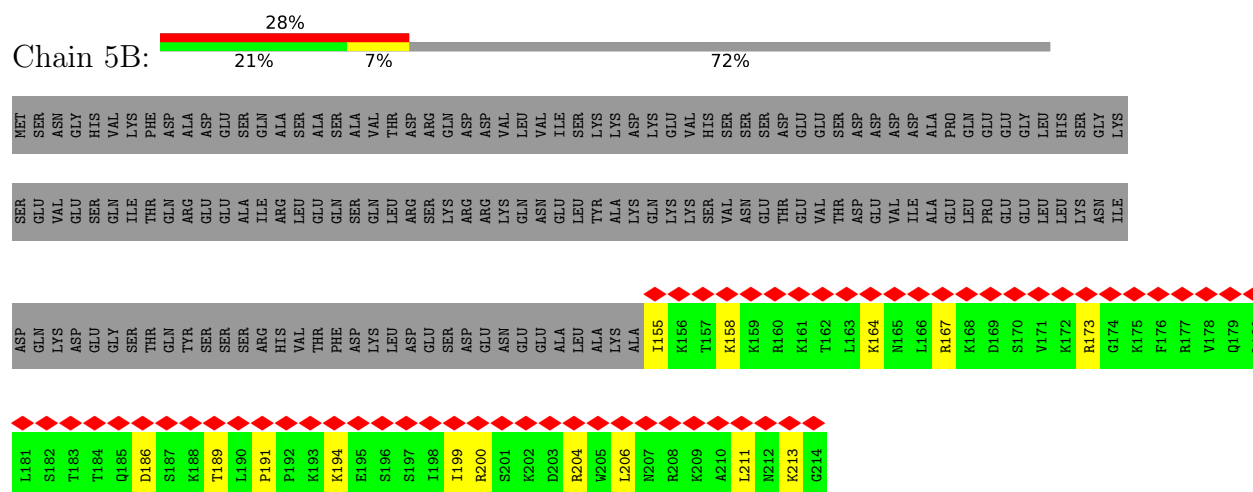
- Molecule 37: U3 small nucleolar RNA-associated protein 6

Chain B6:



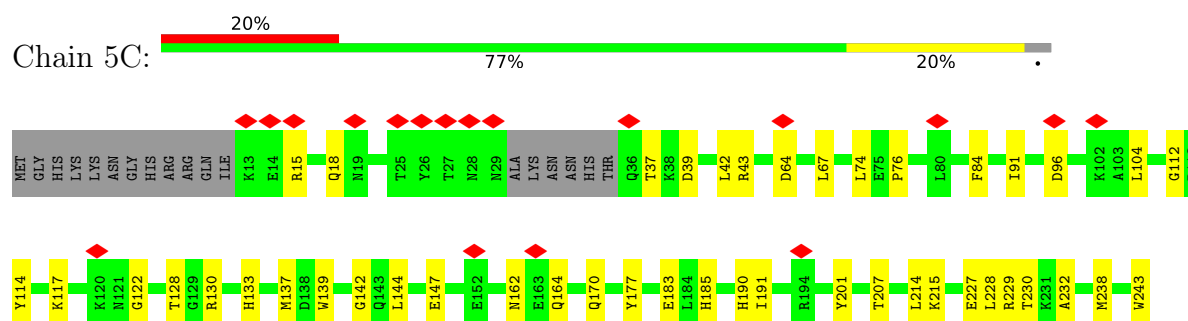
- Molecule 38: Bud site selection protein 21

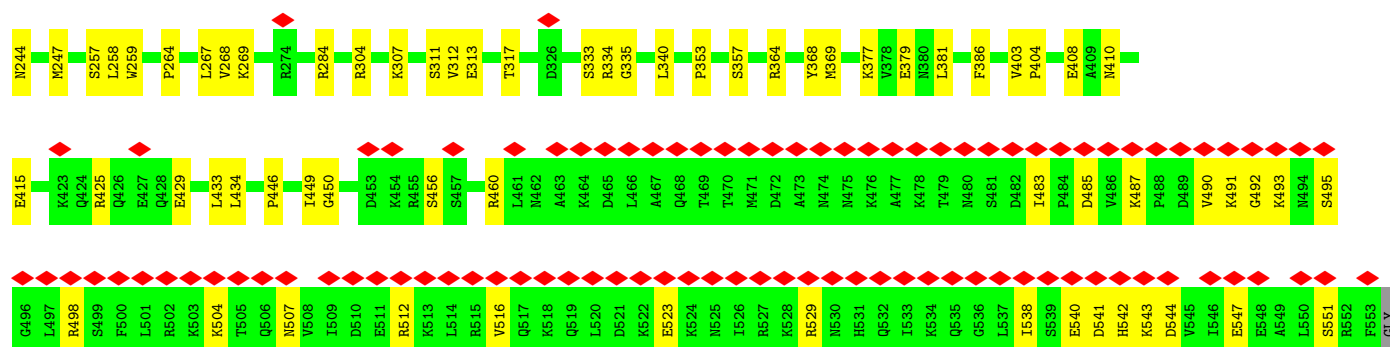
Chain 5B:



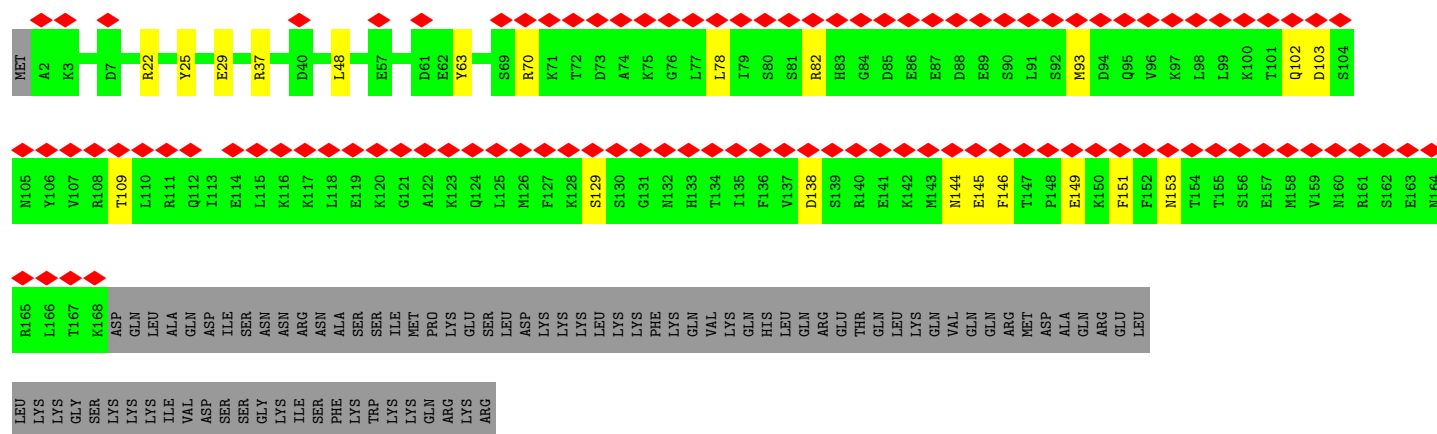
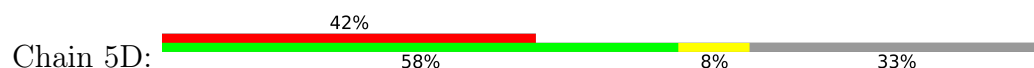
- Molecule 39: U3 small nucleolar RNA-associated protein 7

Chain 5C:

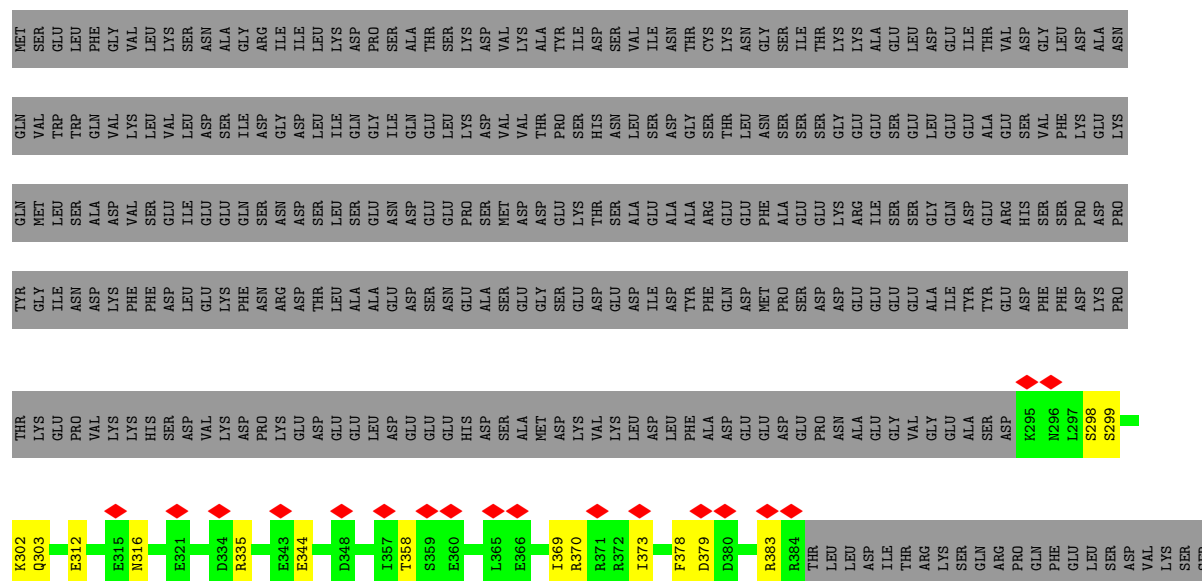




• Molecule 40: U3 small nucleolar RNA-associated protein 11

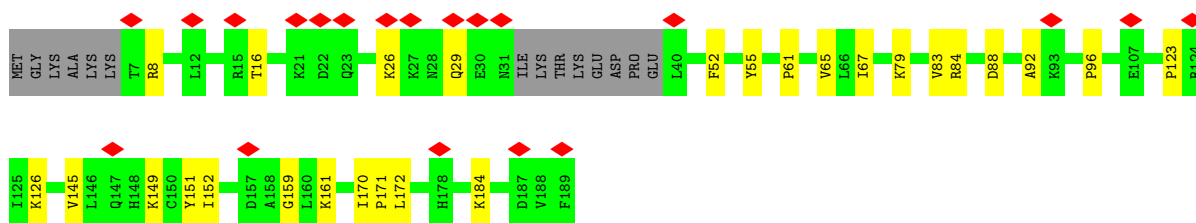


• Molecule 41: U3 small nucleolar RNA-associated protein MPP10

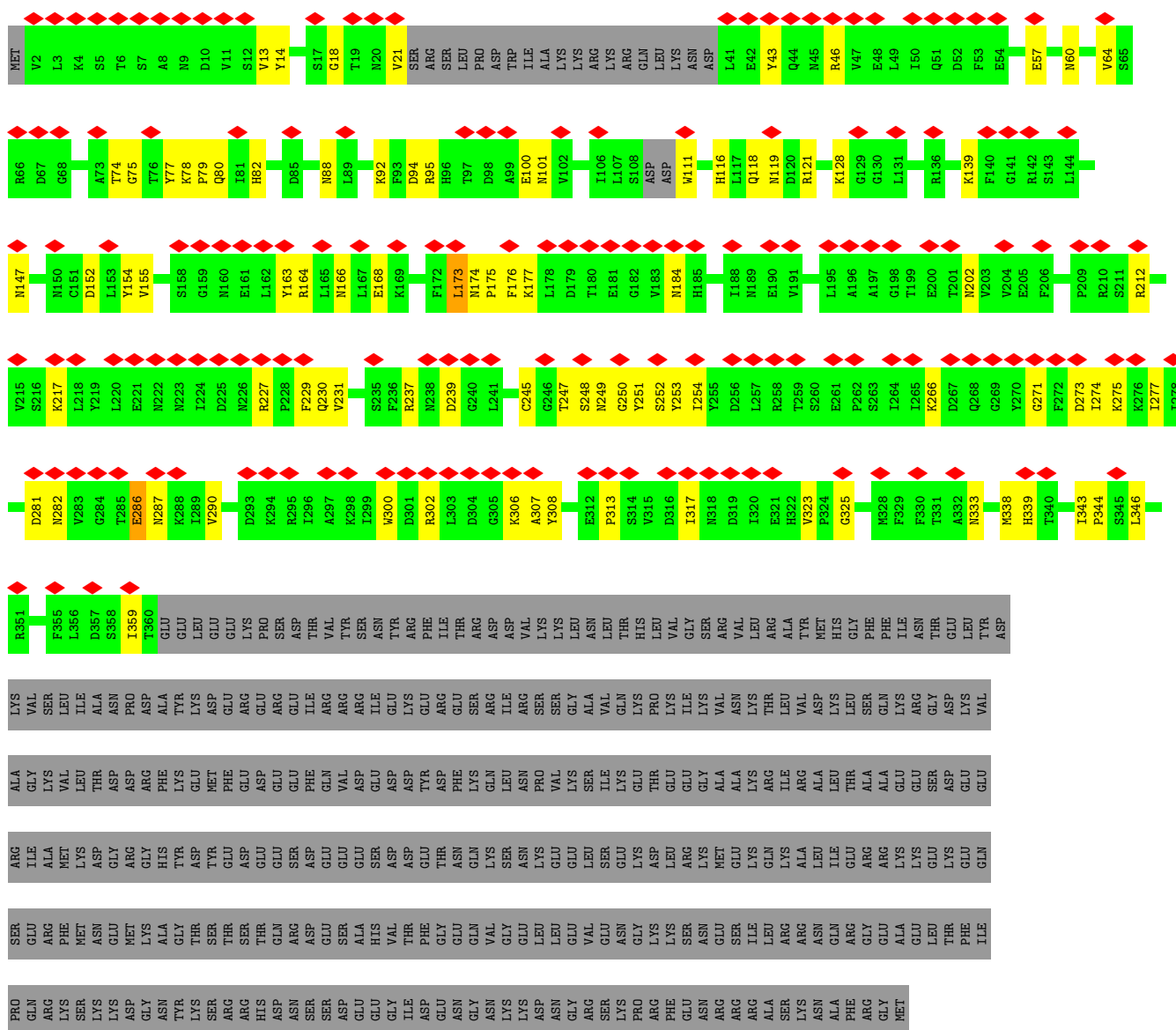




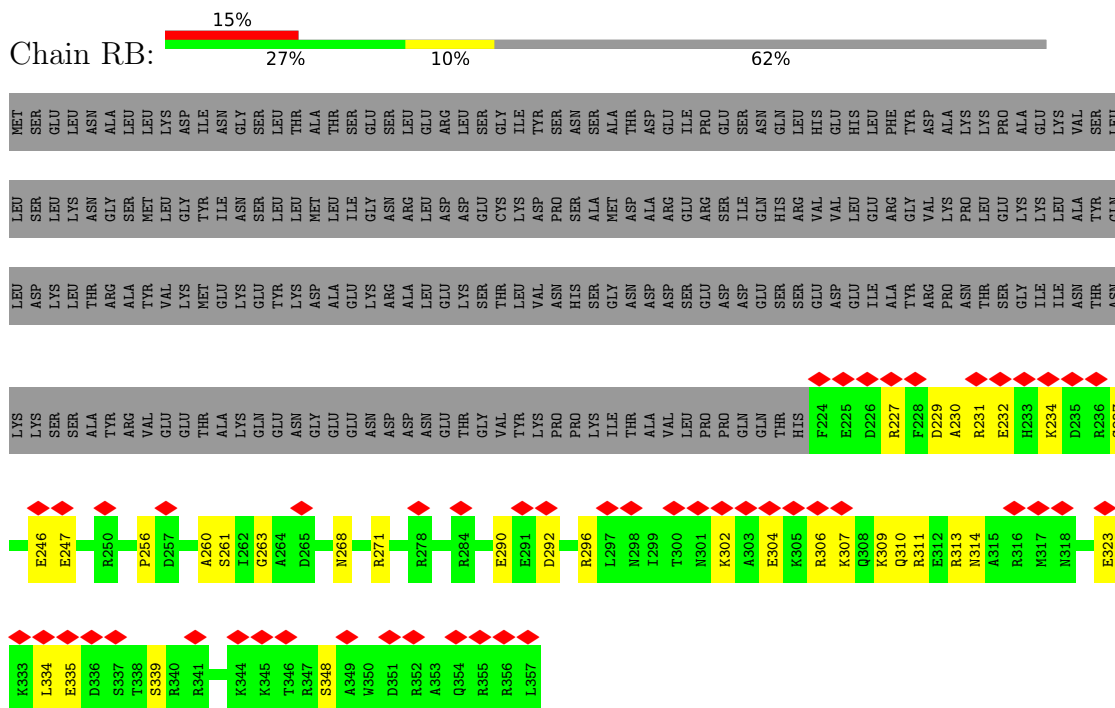
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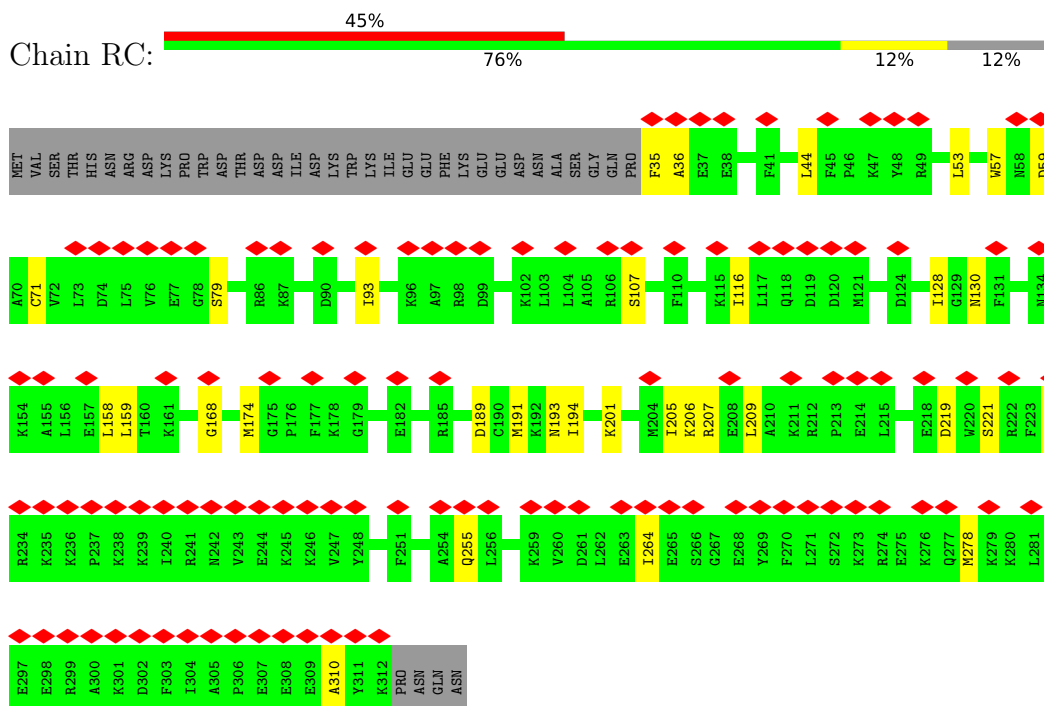
Chain RA:



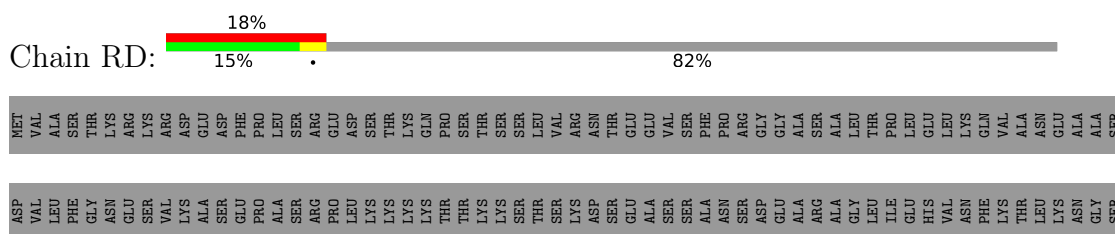
- Molecule 49: U3 small nucleolar ribonucleoprotein protein LCP5



- Molecule 50: KRR1 small subunit processome component

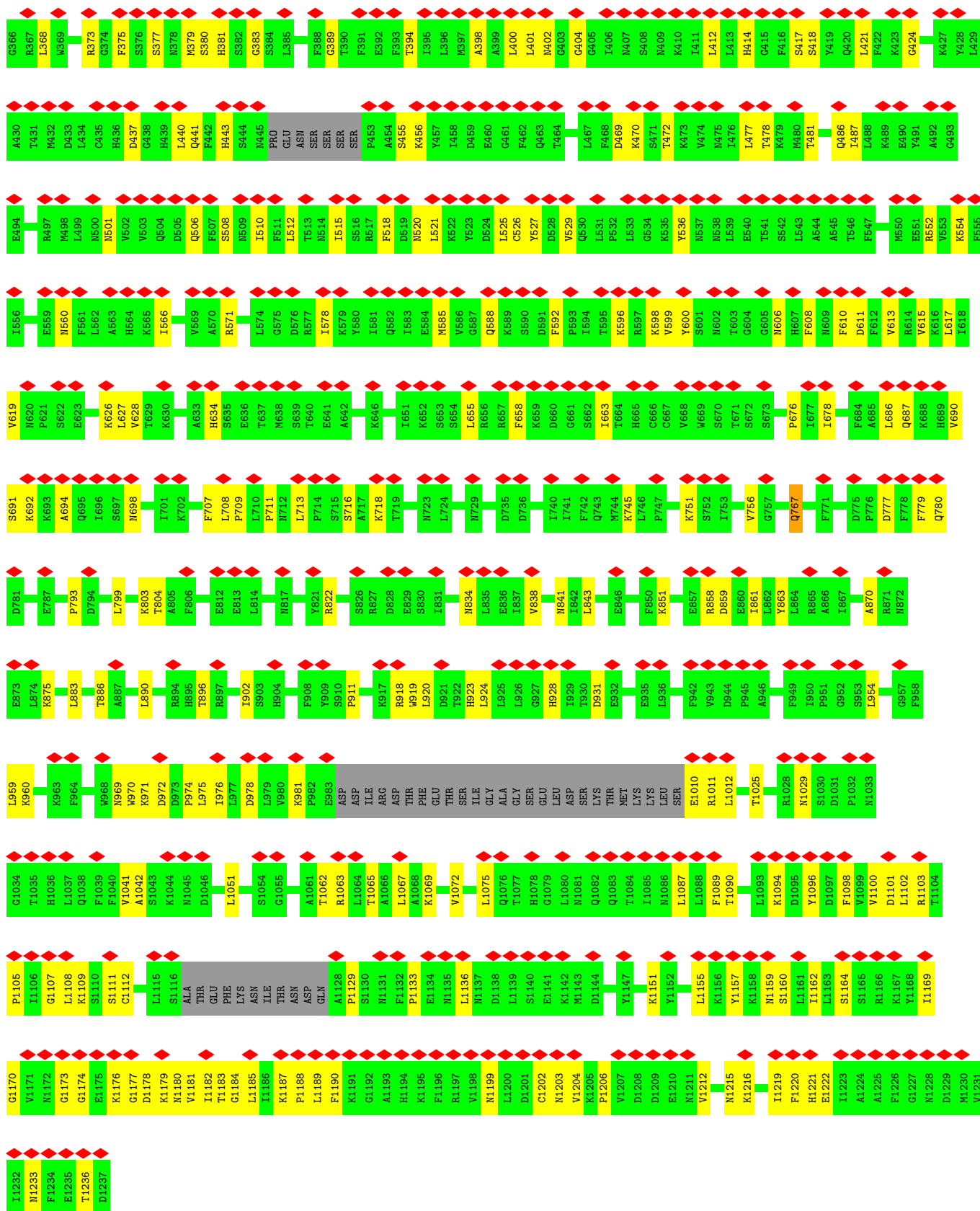


- Molecule 51: rRNA biogenesis protein RRP5

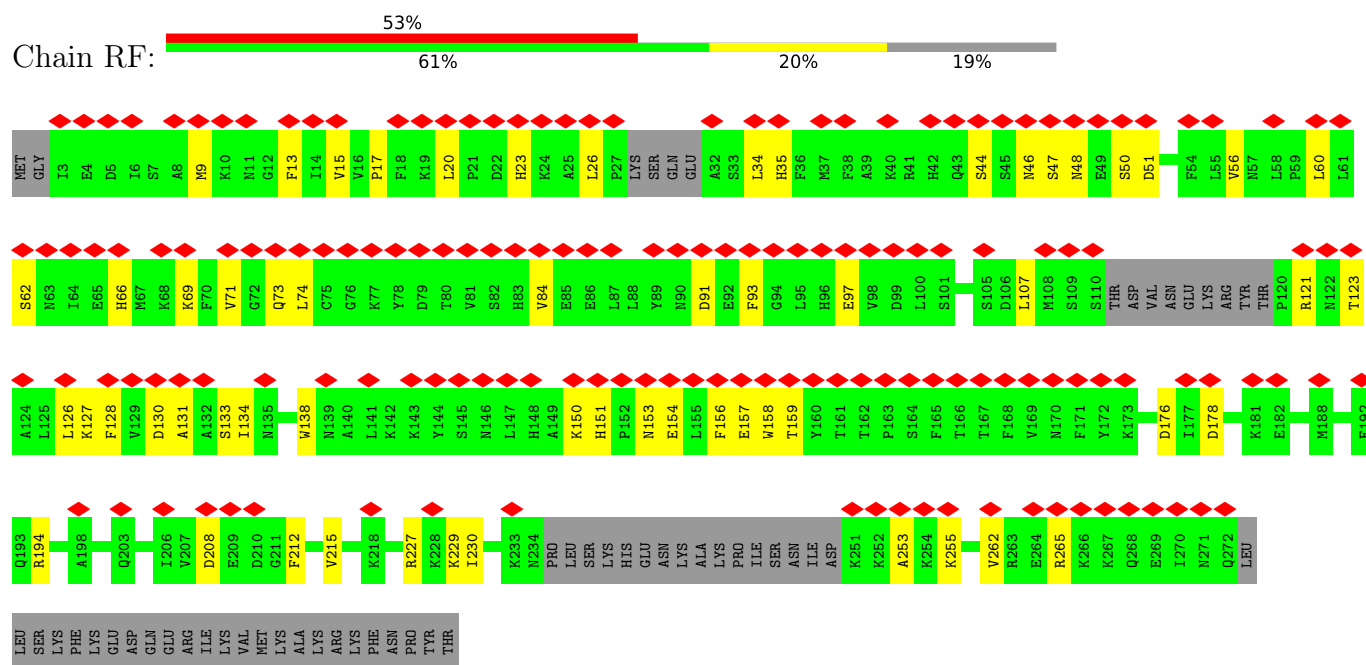




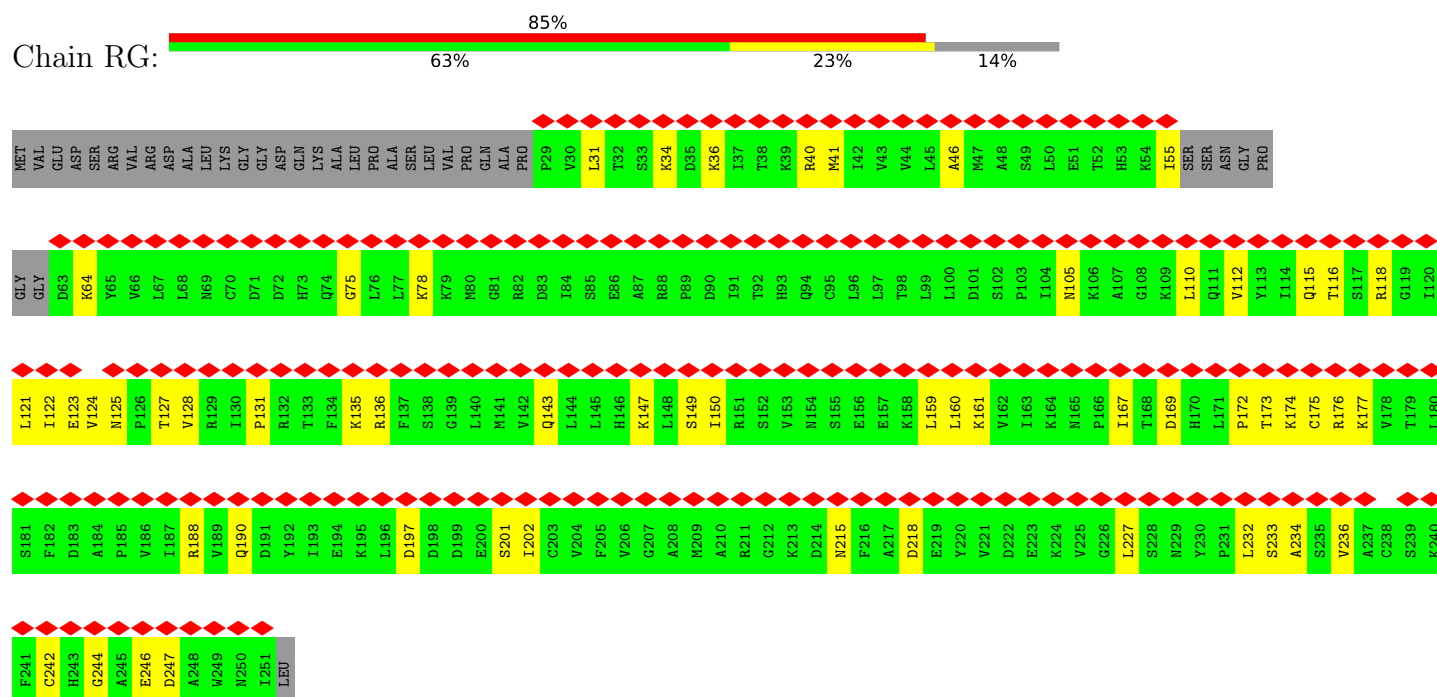




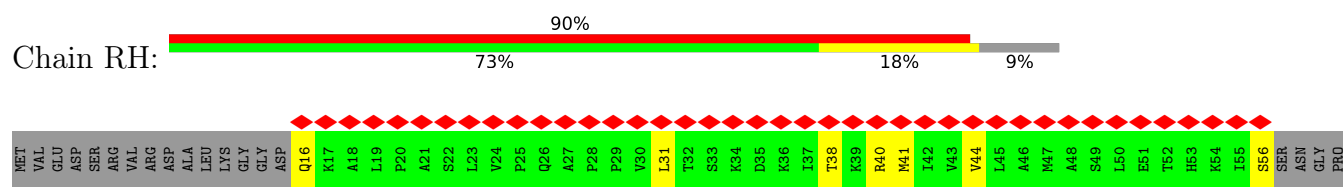
• Molecule 53: Ribosomal RNA-processing protein 7



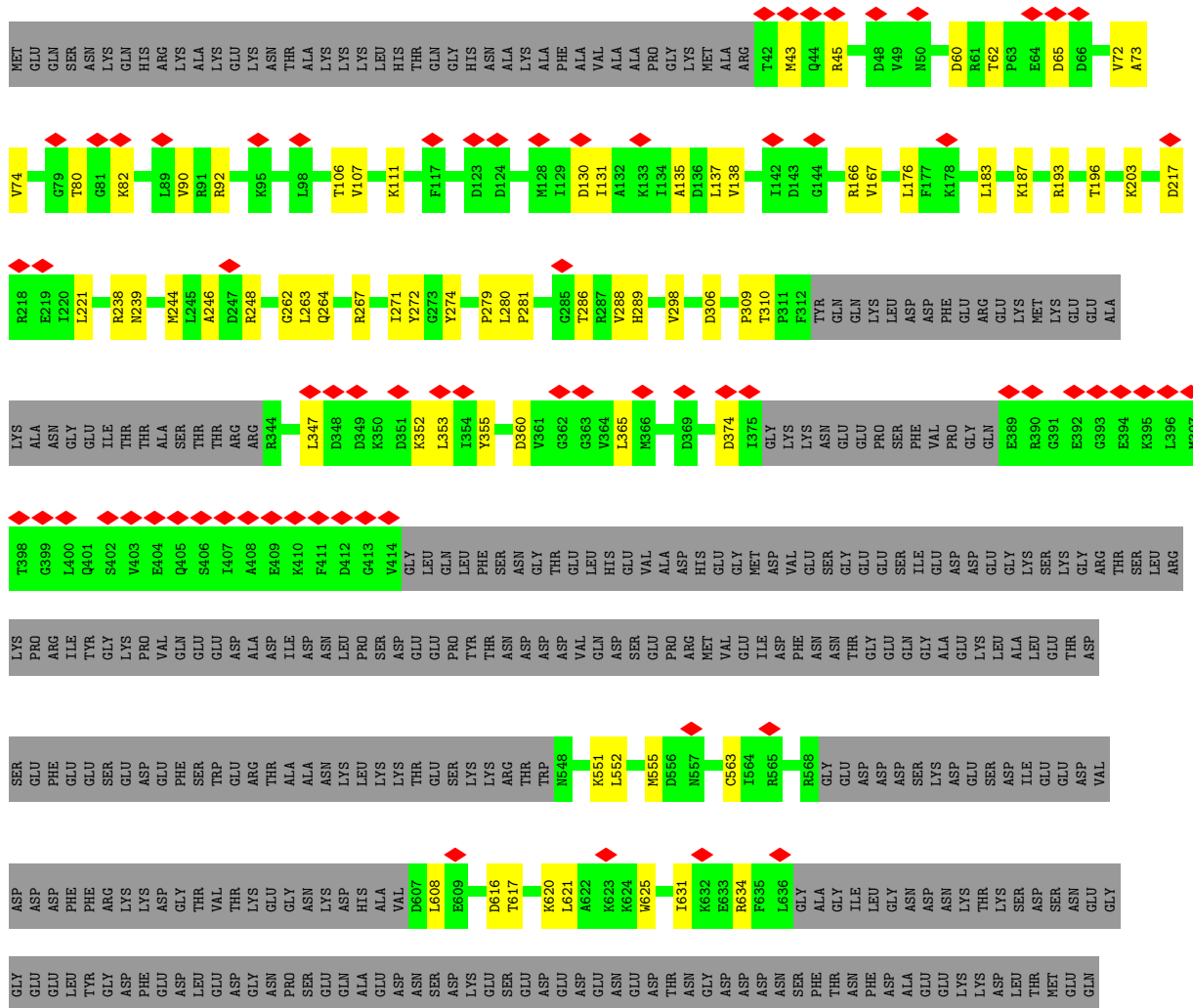
• Molecule 54: Ribosomal RNA small subunit methyltransferase NEP1

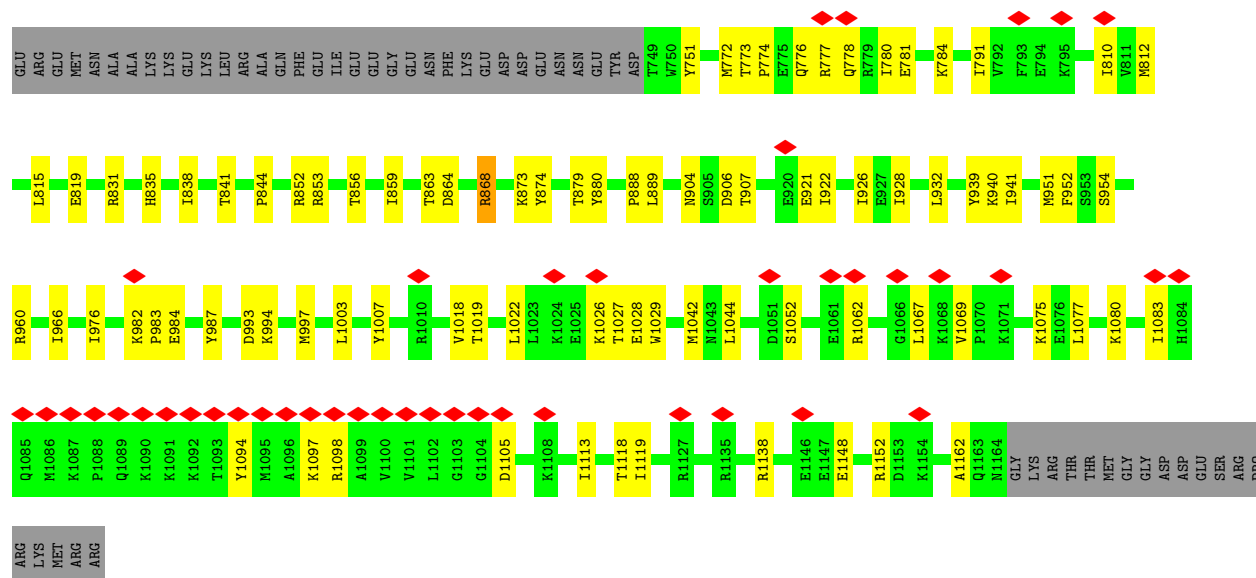


• Molecule 54: Ribosomal RNA small subunit methyltransferase NEP1

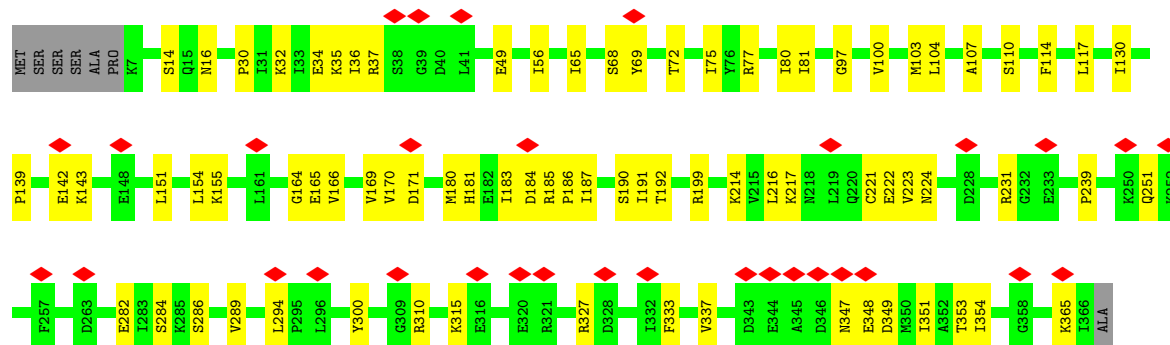
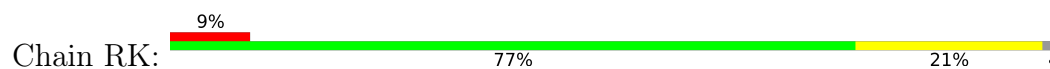


- Molecule 55: Ribosome biogenesis protein BMS1

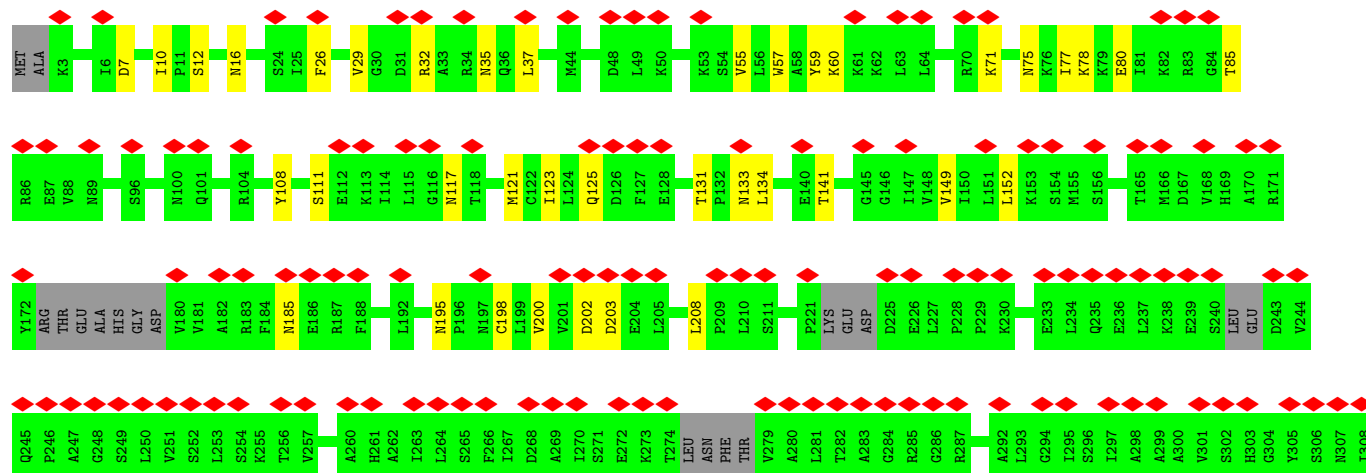
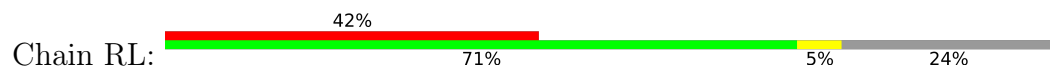


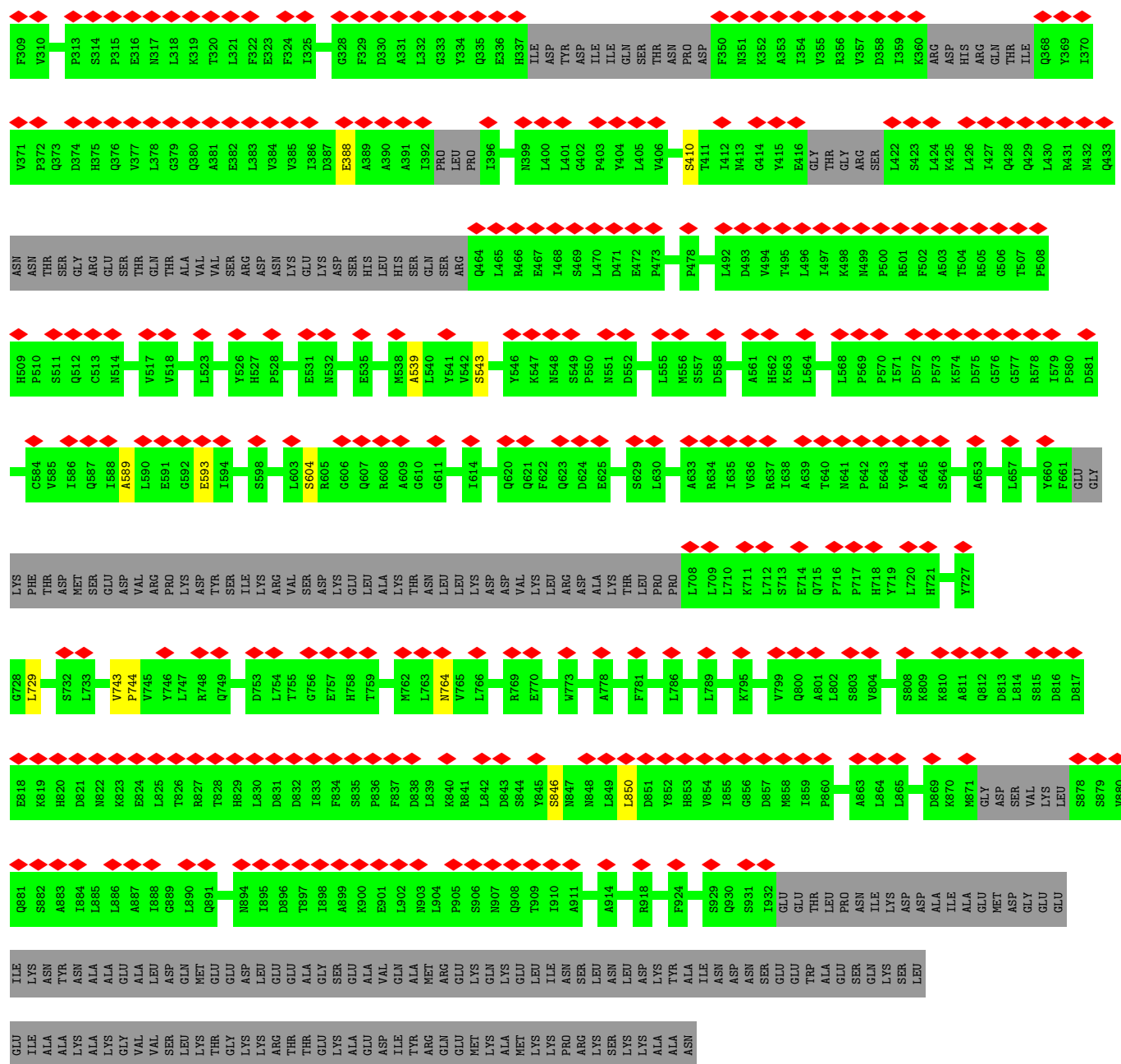


• Molecule 56: RNA 3'-terminal phosphate cyclase-like protein

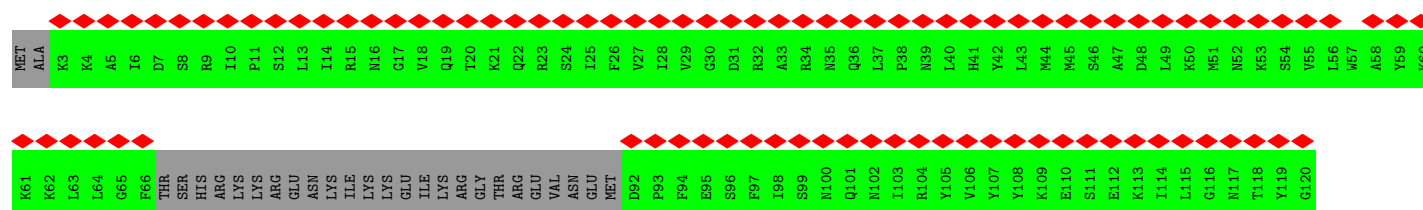
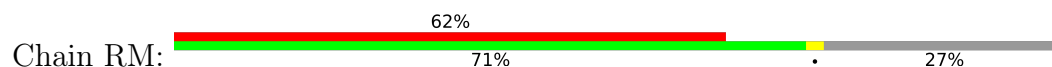


• Molecule 57: RNA cytidine acetyltransferase





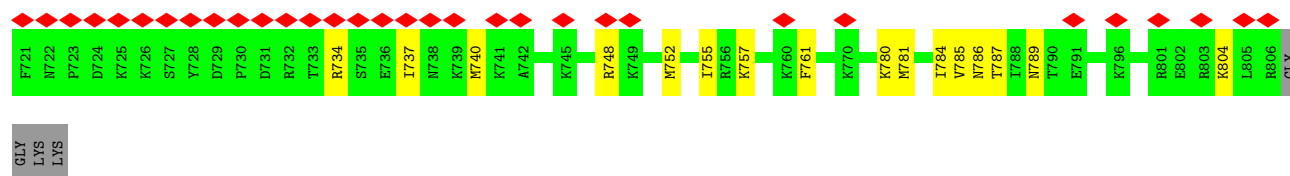
• Molecule 57: RNA cytidine acetyltransferase



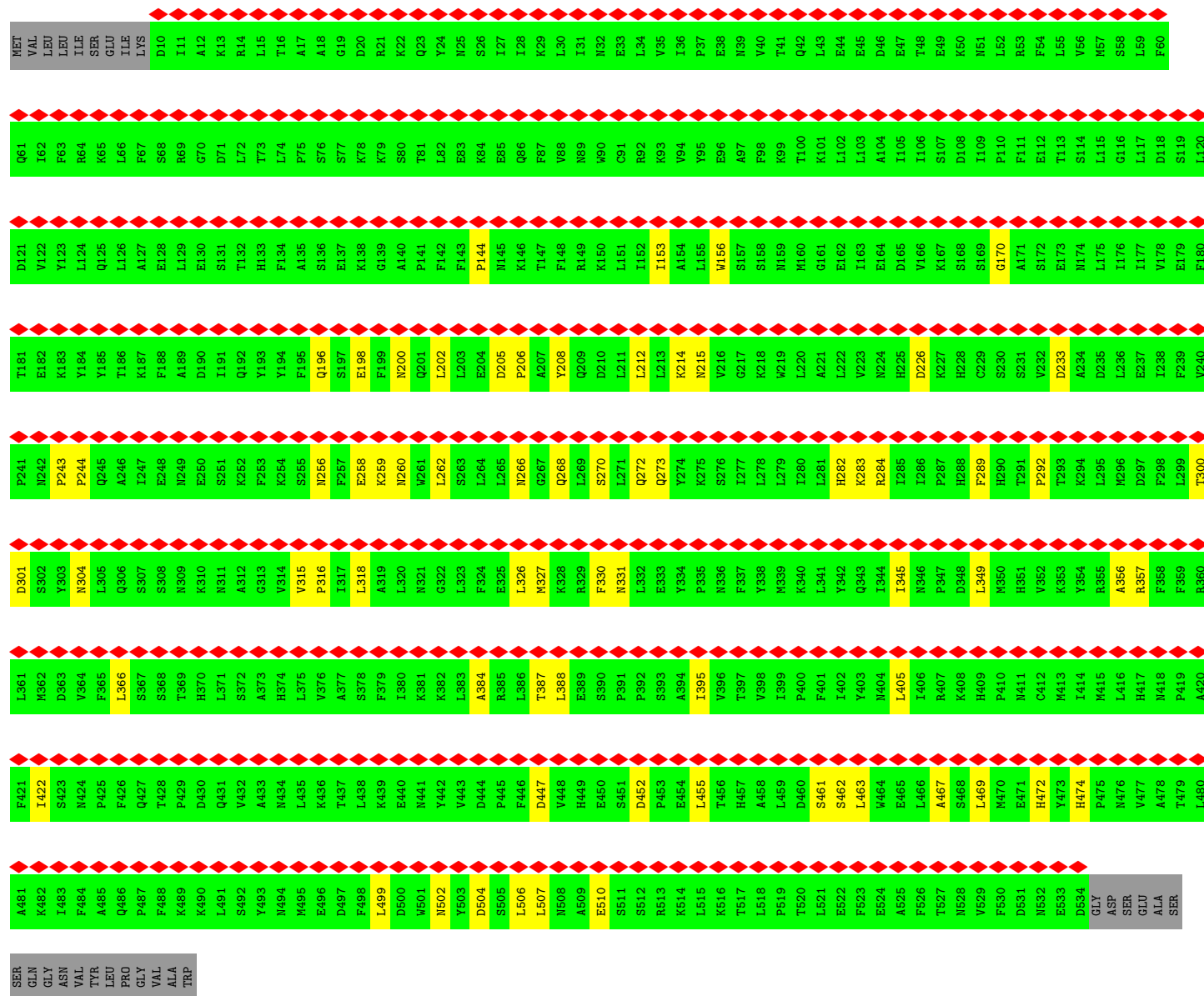
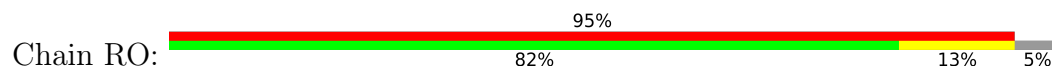


Chain RN:

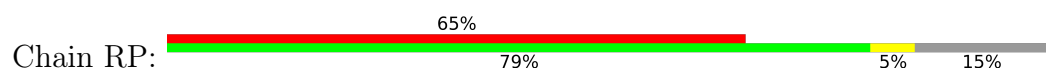




• Molecule 59: Nucleolar complex protein 4



• Molecule 60: U3 small nucleolar RNA-associated protein 20



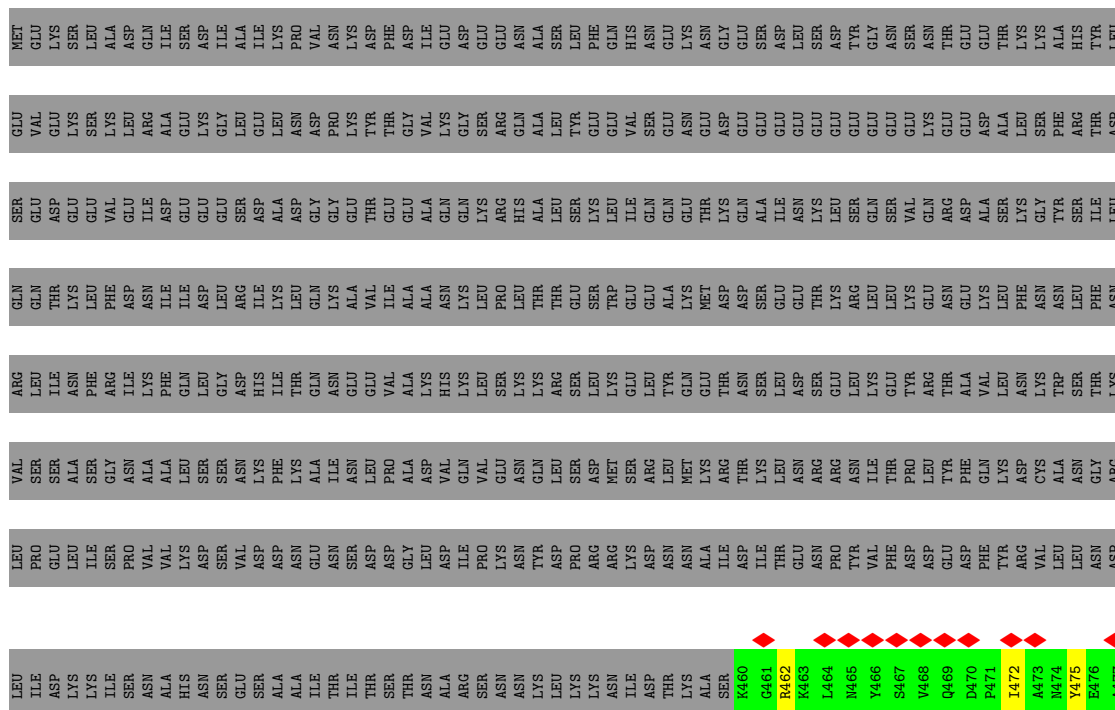
MET	ALA	LYS	GLN	ARG	GLN	THR	THR	LYS	SER	SER	LYS	ARG	TYR	ARG	TYR	SER	S18	F19	K20	A21	K22	I23	D24	D25	L26	K27	T28	E29	P30	A31	ARG	ASN	LEU	GLU	LYS	ARG	VAL	HIS	ASP	TYR	VAL	GLU	SER	S45	H46	F47	L48	A49	S50	F51	D52	Q53	W54	A55	E56	L57	N58	L59	S60																																						
A61	K62	P63	T64	E65	F66	A67	A68	E69	L70	E71	H72	T131	H73	L132	V74	T75	L77	P78	Q79	I80	L81	H82	H83	D84	K85	K86	I87	F88	N89	S90	L91	V92	S93	P94	I95	N96	I97	H98	D99	E100	F101	S102	L103	Q104	P105	L106	L107	D108	L109	L110	A111	Q112	F113	C114	H115	D116	L117	G118	P119	D120																																					
F121	L122	K123	F124	Y125	E126	E127	A128	I129	F131	K130	T131	H132	L133	N134	L135	L136	D137	A138	I139	I140	E141	F142	E143	S144	S145	N146	V147	F148	E149	V150	G151	F152	L155	A156	Y157	I158	F159	K160	Y161	L162	S163	K164	F165	L166	V167	K168	L169	L170	V171	L172	T173	C174	D175	L176	L177	I178	P179	L180																																							
S182	H183	S184	K185	E186	Y187	L188	S189	R190	F191	M255	S192	A193	E194	A195	L196	S197	F198	L199	V200	K201	K202	C203	P204	V205	S206	H207	S273	R209	E210	R213	S214	V215	F216	E217	I282	K218	L219	E220	GLY	ASP	ASP	GLU	GLN	THR	ASN	L228	Y229	E230	G231	L232	L233	I234	T237	E238	S239	M240	T241	E245																																							
T246	L247	H248	S249	K250	A251	K252	A253	I254	M255	S256	V257	L258	L259	H260	E261	A262	L263	T264	K265	S266	P268	E269	R270	S271	V272	S273	L274	D277	L278	W279	M280	N281	E282	I283	S283	K284	Y285	A286	S287	I288	E289	S290	L291	K292	P293	V294	Y295	E296	V297	M298	Y299	Q300	D301	F302	N303	D304	S305	LEU																																							
ASP	ALA	THR	ASN	ASP	R313	I314	L315	K316	V317	L318	T319	T320	I321	V322	F323	S324	E325	SER	GLY	ARG	LYS	I330	P331	D332	W333	N334	K335	I336	T337	I338	L339	I340	F341	R342	I343	M344	SER	GLN	GLU	ASN	CYS	ALA	SER	L353	S354	Q355	D356	K357	V358	A359	F360	L361	F362	A363	L364	F365	I366																																								
R367	N368	S369	D370	V371	K372	T373	T375	L376	F377	H378	Q379	K380	L381	F382	N383	Y384	A385	L386	T387	N388	S389	S390	D391	W392	C392	F393	L394	E395	F396	F397	Q398	F399	A400	L401	R402	L403	S404	Y405	E406	R407	V408	F409	S410	F411	ASN	GLY	LEU	L419	F420	L421	K422	M423	M424	W425	Q426																																										
S427	Q428	Q429	K430	K431	I432	A433	L434	F435	F436	L437	E438	V439	D440	D441	K442	P443	E444	L445	Q446	K447	W448	R449	E450	W451	M452	F453	P454	E455	E456	F457	I458	L459	S460	I461	R462	D463	F464	F465	V466	T467	A468	E469	I470	N471	D472	S473	M474	D475	L476	F477	E478	I479	TYR	TRP	ARG	ALA	ILE	I485	F486																																						
K487	Y488	S489	K490	L491	Q492	N493	T494	E495	I496	I497	ILE	PRO	LEU	LEU	E502	R503	I504	T507	F508	A509	S510	P511	D512	N513	F514	T515	K516	ASP	VAL	P579	G520	T521	L522	L523	K524	R527	K528	E529	D530	D531	A532	S533	ASN	ASN	LEU	L538	K539	T540	I541	L542	D543	N544	Y545	E546	N547	Y548																																									
K549	E550	S551	L552	ASN	PHE	LEU	ARG	G557	W558	N559	K560	L561	V562	S563	N564	L565	H566	P567	S568	E569	S570	L571	K572	G573	L574	M575	S576	H577	Y578	P579	S580	L581	L582	L583	S584	L585	T586	D587	N588	F589	M590	L591	P592	D593	G594	K595	I596	R597	GLU	THR	LEU	GLU	MET	K605	T606	L607	M608																																								
I609	L610	Q611	G612	M613	Q614	V615	P616	D617	L618	L619	S620	S621	C622	M623	V624	I625	H626	E627	Q633	N634	A635	T639	I640	R641	I642	K643	ASN	VAL	GLY	ALA	GLU	PHE	GLY	L682	L683	L684	L685	L686	G687	G688	L689	L690	L691	L692	L693	L694	L695	L696	L697	L698	L699	L700	L701	L702	L703	L704	L705	L706	L707	L708	L709	L710	L711	L712	L713	L714	L715	L716	L717	L718	L719	L720	L721	L722	L723	L724	L725	L726	L727	L728	L729	L730	L731	L732	L733	L734	L735	L736	L737	L738	L739	L740	L741	L742	L743	L744	L745
I746	W747	S748	K749	Y750	S751	T752	Q753	N754	T755	S756	I757	F758	S759	T760	T761	I762	E763	R764	R765	G766	T767	T768	T769	Y770	T771	I772	L773	I774	R775	N776	Q777	A778	L779	K780	W781	M782	I785	P786	A789	E790	N791	H792	F793	V794	A797	P798	F799	V800	Y801	N802	D803	K808	E811	D812																																											
M813	E814	N815	E816	I819	T820	G821	S822	L832	K833	T834	L835	S836	K837	F838	I841	V844	Y845	S846	A847	T848	E849	L850	H851	D852	H853	L854	L857	L858	R861	N862	T863	D864	V865	Q866	K867	L868	A869	L870	D871	A872	L873	L874	A875	Y876	K877	N878	P879	T880	L881	N882	K883	Y884	R885																																												

D886	N887	L888	K889	D893	L896	D899	E900	I901	T902	T903	F904	L905	T906	N908	G909	S910	Q911	S912	I913	K914	A915	E916	D917	E918	K919	V920	M922	P923	Y924	I928	R932	A933	Q934	V935	F936	P937	T938	S939	G940	Q941	K942	R943	S944	R945	K946	I947	A948	V949	I950	S951	V952	L953						
P954	N955	I961	F964	R971	I974	N975	I976	F977	F978	G979	S986	A989	T990	T993	I994	R995	T998	V1001	M1002	I1003	V1004	M1005	S1006	T1007	H1018	T1019	V1022	P1025	L1026	S1029	I1030	A1031	M1032	A1033	Y1034	Y1035	V1036	L1037	E1040	S1041	T1042	E1043	E1044	V1045	H1046													
L1047	R1048	K1049	M1050	A1051	R1055	Q1056	Q1057	K1060	C1061	L1062	S1063	S1064	V1065	F1066	F1068	F1073	D1074	W1075	S1076	T1077	S1078	Y1083	A1084	V1085	V1086	V1087	K1088	P1089	I1091	S1092	H1093	F1094	S1095	D1096	E1097	N1098	L1099	Q1100	S1103	S1104	L1105	L1108	F1109	L1110	Y1111	W1112	A1113	H1114	N1115	P1116	S1117							
L1118	Y1119	F1121	M1122	Y1123	Y1124	F1127	A1128	T1129	A1130	T1131	A1132	L1133	M1134	D1135	T1136	I1137	S1138	M1139	Q1140	H1141	V1142	K1143	E1144	A1145	V1146	I1147	G1148	P1149	I1150	I1151	E1152	A1153	A1154	D1155	S1156	I1157	I1158	R1159	M1160	P1161	V1162	M1163	D1164	D1165	H1166	Y1167	V1168	L1169	L1170	V1171	H1172	L1173	I1174	C1175	S1176	C1178		
L1179	K1180	I1181	L1182	P1183	S1184	L1185	Y1186	V1187	L1188	S1189	D1191	S1192	M1193	S1194	I1195	S1196	T1197	F1198	L1199	N1200	L1201	L1202	V1203	S1204	I1205	T1206	E1207	M1208	G1209	F1210	I1211	Q1212	D1213	D1214	H1215	V1216	R1217	S1218	R1219	L1220	I1221	S1222	S1223	L1224	I1225	S1226	I1227	L1228	K1229	G1230	K1231	L1232	K1233	K1234	L1235	Q1236	E1237	N1238
D1239	T1240	Q1241	K1242	L1243	L1244	K1245	L1246	L1247	K1248	L1249	I1250	V1251	F1252	N1253	Y1254	N1255	C1256	S1257	W1258	S1259	D1260	I1261	E1262	E1263	L1264	Y1265	T1266	T1267	I1268	S1269	L1270	L1271	F1272	K1273	T1274	F1275	D1276	E1277	R1278	N1279	L1280	R1281	V1282	S1283	L1284	T1285	E1286	L1287	F1288	I1289	E1290	L1291	G1292	R1293	K1294	E1297	L1298	E1299
S1300	T1301	S1302	K1303	L1304	V1305	A1306	D1307	L1308	N1309	S1310	Y1311	S1312	S1313	S1314	R1315	W1316	H1317	E1318	Y1319	D1320	F1321	P1322	ARG	I1323	LEU	SER	PHE	LYS	GLY	LEU	I1328	GLU	ASP	GLY	TYR	SER	GLU	LEU	TRP	PRO	LEU	PHE	THR	PHE	LEU	HIS	PHE	ILE	ASN	LYS	GLN	GLU						
GLU	LEU	ALA	LEU	THR	ASN	ALA	SER	HIS	ALA	ILE	MET	LYS	PHE	I1376	F1377	I1378	N1379	E1380	K1381	P1382	N1383	L1384	N1385	I1386	A1387	S1388	K1389	S1390	I1391	S1392	M1393	L1394	K1395	D1396	I1397	L1398	L1399	P1400	N1401	I1402	R1403	I1404	G1405	L1406	R1407	D1408	S1409	L1410	E1411	E1412	V1413	GLN	SER	GLU	THR	VAL	SER	
V1420	L1421	S1422	Y1423	M1424	V1425	K1426	N1427	T1428	K1429	Y1430	F1431	T1432	D1433	D1436	I1439	L1440	L1441	Y1442	N1443	G1444	D1445	E1446	E1447	A1448	D1449	PHE	PHE	THR	ASN	VAL	ASN	ASN	HIS	ILE	GLN	LEU	HIS	R1461	R1462	Q1463	R1464	A1465	I1466	K1467	E1471	H1472	A1473	L1474	Q1475	L1476	K1477	D1478	M1479	S1480	T1481	L1482	H1483	
Y1484	L1485	I1486	M1488	I1489	E1490	H1491	Y1492	V1493	F1494	S1495	D1496	D1497	E1498	R1499	Y1500	R1501	N1502	I1503	G1504	E1506	T1507	Q1508	I1509	A1510	I1511	G1512	G1513	Q1516	H1517	M1518	S1519	V1520	N1521	Q1522	Y1523	K1524	A1525	L1526	L1527	R1528	R1529	Y1530	I1531	S1532	M1533	L1534	K1535	T1536	K1537	P1538	M1539	Q1540	M1541	K1542	Q1543	A1544		
V1545	Q1546	L1547	I1548	VAL	GLN	LEU	SER	VAL	PRO	LEU	ARG	GLU	T1558	L1559	R1560	I1561	V1562	R1563	D1564	G1565	A1566	E1567	S1568	K1569	L1570	S1573	K1574	F1575	P1576	S1577	ASN	LEU	ASP	GLU	PRO	SER	ASN	PHE	ILE	LYS	GLN	GLU	LEU	TYR	PRO	THR	SER	LYS	ILE	LEU	GLY	THR	ARG	ASP	GLU	THR		
I1571	I1572	I1573	I1574	I1575	I1576	I1577	C1640	Q1641	V1642	L1643	R1644	S1645	K1646	S1647	E1648	E1649	L1650	R1651	D1652	V1653	V1654	R1655	Y1656	I1663	T1664	L1665	E1668	V1669	L1670	V1671	F1672	H1673	T1674	K1675																								
E1676	A1679	T1680	L1681	K1682	R1683	G1684	S1685	Q1686	I1687	H1688	V1689	L1690	S1691	Y1692	H1695	Y1696	M1701	H1702	G1703	V1704	L1705	P1706	H1707	S1708	D1709	L1710	D1711	T1712	S1713	M1716	K1719	I1720	I1721	H1722	E1723	N1724	I1725	F1726	G1727	F1728	A1729	G1730	E1731	E1732	K1733	D1734	S1735	E1736	N1737	Y1738	H1739	T1740	K1741	V1742				

- Molecule 61: U3 small nucleolar RNA-associated protein 14



- Molecule 65: Protein BFR2



- Molecule 66: Unassigned helices





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	17924	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.126	Depositor
Minimum map value	-0.059	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	597.632, 597.632, 597.632	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.334, 1.334, 1.334	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	3A	0.45	0/4141	0.48	0/6433
2	5A	0.44	0/8952	0.47	1/13931 (0.0%)
3	SA	0.35	0/31535	0.47	5/49095 (0.0%)
4	SC	0.44	0/1856	0.77	3/2490 (0.1%)
5	SF	0.29	0/1854	0.66	0/2504
6	SG	0.49	0/1690	0.67	2/2285 (0.1%)
7	SH	0.24	0/1341	0.61	1/1789 (0.1%)
8	SI	0.32	0/1341	0.69	2/1806 (0.1%)
9	SJ	0.26	0/1347	0.60	0/1801
10	SK	0.44	0/1410	0.63	0/1888
11	SM	0.26	0/1020	0.64	0/1374
12	SO	0.39	0/1109	0.63	0/1495
13	SP	0.46	0/879	0.79	0/1186
14	SR	0.52	0/990	0.77	3/1335 (0.2%)
15	SX	0.46	0/1020	0.63	0/1371
16	SY	0.48	0/798	0.65	0/1065
17	SZ	0.36	0/822	0.73	0/1103
18	Sc	0.40	0/613	0.63	0/828
19	Sd	0.50	0/499	0.72	2/670 (0.3%)
20	3B	0.51	0/1901	0.66	1/2567 (0.0%)
20	3C	0.37	0/1796	0.66	0/2424
21	3D	0.40	0/2891	0.62	0/3895
22	3E	0.38	0/3059	0.62	0/4153
23	3F	0.36	0/3715	0.64	0/5001
24	3G	0.51	0/928	0.72	1/1262 (0.1%)
24	3H	0.45	0/928	0.70	0/1262
25	A4	0.43	0/5321	0.66	0/7207
26	A5	0.44	0/4044	0.65	0/5493
27	A8	0.23	0/3328	0.62	0/4565
28	A9	0.28	0/951	0.60	0/1287
29	AE	0.32	1/10049 (0.0%)	0.56	1/13737 (0.0%)
30	AF	0.48	0/3993	0.68	0/5413

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
31	AG	0.42	0/6699	0.64	3/9077 (0.0%)
32	B1	0.58	0/6780	0.69	0/9175
33	B2	0.39	0/6853	0.68	2/9256 (0.0%)
34	B3	0.35	0/6014	0.70	2/8137 (0.0%)
35	B8	0.55	0/3848	0.63	0/5218
36	BE	0.53	0/6948	0.64	2/9391 (0.0%)
37	B6	0.37	0/2849	0.60	0/3853
38	5B	0.31	0/499	0.62	0/659
39	5C	0.55	0/4321	0.66	2/5832 (0.0%)
40	5D	0.47	0/1417	0.73	0/1885
41	5E	0.44	0/1665	0.66	2/2233 (0.1%)
42	5F	0.64	0/1559	0.72	1/2097 (0.0%)
43	5G	0.52	0/2337	0.67	2/3148 (0.1%)
44	5H	0.47	0/601	0.59	0/789
45	5I	0.56	0/3844	0.68	0/5174
46	5J	0.36	0/1302	0.58	0/1728
47	5K	0.51	0/1426	0.66	0/1917
48	RA	0.27	0/2769	0.66	2/3753 (0.1%)
49	RB	0.37	0/1121	0.79	0/1487
50	RC	0.40	0/2245	0.60	0/3021
51	RD	0.26	0/2453	0.67	2/3308 (0.1%)
52	RE	0.33	0/8924	0.64	4/12070 (0.0%)
53	RF	0.28	0/2004	0.66	1/2697 (0.0%)
54	RG	0.32	0/1727	0.70	0/2329
54	RH	0.37	0/1828	0.59	0/2470
55	RJ	0.44	0/6514	0.61	0/8768
56	RK	0.39	0/2832	0.64	0/3825
57	RL	0.21	0/4549	0.51	1/6241 (0.0%)
57	RM	0.15	0/3765	0.44	1/5218 (0.0%)
58	RN	0.30	0/4591	0.63	1/6187 (0.0%)
59	RO	0.32	0/3849	0.63	0/5261
60	RP	0.21	0/12225	0.54	4/16812 (0.0%)
61	RQ	0.41	0/1678	0.63	1/2282 (0.0%)
62	RS	0.30	0/2104	0.70	0/2854
63	RT	0.38	0/1379	0.68	0/1853
64	RV	0.43	0/1456	0.60	1/1937 (0.1%)
65	RY	0.23	0/307	0.53	0/415
All	All	0.40	1/233403 (0.0%)	0.60	56/325072 (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
4	SC	0	1
5	SF	0	2
8	SI	0	3
9	SJ	0	1
11	SM	0	1
12	SO	0	1
13	SP	0	1
17	SZ	0	1
18	Sc	0	1
21	3D	0	3
22	3E	0	1
23	3F	0	1
24	3G	0	2
24	3H	0	1
25	A4	0	1
26	A5	0	1
27	A8	0	4
31	AG	0	2
32	B1	0	3
33	B2	0	9
34	B3	0	11
36	BE	0	1
39	5C	0	2
40	5D	0	1
41	5E	0	1
42	5F	0	1
43	5G	0	1
45	5I	0	2
48	RA	0	2
49	RB	0	1
52	RE	0	1
53	RF	0	1
55	RJ	0	2
56	RK	0	1
57	RL	0	1
57	RM	0	1
58	RN	0	1
59	RO	0	1
60	RP	0	3
61	RQ	0	1
64	RV	0	2
All	All	0	78

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
29	AE	564	VAL	C-N	5.39	1.40	1.34

The worst 5 of 56 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	RD	1223	PRO	N-CA-CB	8.04	111.69	103.25
52	RE	1105	PRO	O-C-N	-7.87	112.02	122.64
51	RD	1205	PRO	N-CA-CB	7.68	110.60	103.15
33	B2	267	ASP	CA-C-N	6.86	134.65	121.54
33	B2	267	ASP	C-N-CA	6.86	134.65	121.54

There are no chirality outliers.

5 of 78 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	SC	238	GLU	Peptide
5	SF	193	GLY	Peptide
5	SF	195	ILE	Peptide
8	SI	31	SER	Peptide
8	SI	64	VAL	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	3A	3711	0	1882	21	0
2	5A	8009	0	4030	52	0
3	SA	28210	0	14226	240	0
4	SC	1830	0	1914	33	0
5	SF	1815	0	1870	45	0
6	SG	1669	0	1724	18	0
7	SH	1327	0	1403	30	0
8	SI	1321	0	1387	23	0
9	SJ	1324	0	1344	48	0
10	SK	1388	0	1467	32	0
11	SM	997	0	1048	34	0
12	SO	1087	0	1152	13	0
13	SP	868	0	894	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	SR	973	0	1029	14	0
15	SX	1003	0	1040	16	0
16	SY	786	0	843	8	0
17	SZ	809	0	842	15	0
18	Sc	603	0	623	9	0
19	Sd	497	0	535	3	0
20	3B	1865	0	1910	30	0
20	3C	1763	0	1805	34	0
21	3D	2848	0	2815	46	0
22	3E	3028	0	2813	62	0
23	3F	3643	0	3654	79	0
24	3G	916	0	964	11	0
24	3H	916	0	964	26	0
25	A4	5226	0	5199	96	0
26	A5	3976	0	3919	56	0
27	A8	3307	0	2316	39	0
28	A9	939	0	898	19	0
29	AE	9955	0	7968	103	0
30	AF	3911	0	3906	71	0
31	AG	6570	0	6473	127	0
32	B1	6635	0	6525	103	0
33	B2	6723	0	6698	121	0
34	B3	5919	0	6007	127	0
35	B8	3764	0	3757	58	0
36	BE	6810	0	6787	86	0
37	B6	2800	0	2517	35	0
38	5B	495	0	561	13	0
39	5C	4237	0	4239	85	0
40	5D	1396	0	1407	20	0
41	5E	1647	0	1678	30	0
42	5F	1530	0	1572	30	0
43	5G	2296	0	2325	41	0
44	5H	596	0	661	8	0
45	5I	3765	0	3714	73	0
46	5J	1280	0	1331	23	0
47	5K	1403	0	1484	19	0
48	RA	2709	0	2622	66	0
49	RB	1108	0	1087	27	0
50	RC	2207	0	2229	29	0
51	RD	2412	0	2263	40	0
52	RE	8716	0	8828	240	0
53	RF	1963	0	1942	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
54	RG	1701	0	1767	40	0
54	RH	1799	0	1872	29	0
55	RJ	6379	0	6506	110	0
56	RK	2781	0	2878	51	0
57	RL	4539	0	2874	29	0
57	RM	3779	0	1650	8	0
58	RN	4529	0	4262	69	0
59	RO	3766	0	3269	47	0
60	RP	12171	0	7749	76	0
61	RQ	1651	0	1450	28	0
62	RS	2051	0	2096	53	0
63	RT	1357	0	1426	15	0
64	RV	1448	0	1435	16	0
65	RY	299	0	275	6	0
66	X1	375	0	92	4	0
67	5K	1	0	0	0	0
67	Sc	1	0	0	0	0
68	RJ	32	0	12	2	0
69	RJ	1	0	0	0	0
All	All	226161	0	194704	2951	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 7.

The worst 5 of 2951 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:RE:1203:ASN:HD22	52:RE:1219:ILE:CG2	1.50	1.22
52:RE:1102:LEU:HD23	52:RE:1180:ASN:O	1.40	1.20
52:RE:1216:LYS:NZ	52:RE:1236:THR:HG23	1.65	1.11
52:RE:518:PHE:CZ	52:RE:1075:LEU:HG	1.92	1.05
52:RE:1203:ASN:HD22	52:RE:1219:ILE:HG21	1.18	1.04

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	3A	169/333 (50%)	44 (26%)	2 (1%)
2	5A	365/700 (52%)	104 (28%)	7 (1%)
3	SA	1303/1808 (72%)	494 (37%)	19 (1%)
All	All	1837/2841 (64%)	642 (34%)	28 (1%)

5 of 642 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	3A	2	U
1	3A	14	A
1	3A	15	U
1	3A	24	U
1	3A	25	U

5 of 28 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	SA	372	G
3	SA	1632	C
3	SA	538	A
3	SA	1197	C
3	SA	417	A

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
68	GTP	RJ	1201	69	33,34,34	0.56	0	50,54,54	0.72	1 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
68	GTP	RJ	1201	69	-	3/22/38/38	0/3/3/3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
68	RJ	1201	GTP	C4'-O4'-C1'	-2.15	104.72	109.47

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
68	RJ	1201	GTP	O4'-C4'-C5'-O5'
68	RJ	1201	GTP	C3'-C4'-C5'-O5'
68	RJ	1201	GTP	C4'-C5'-O5'-PA

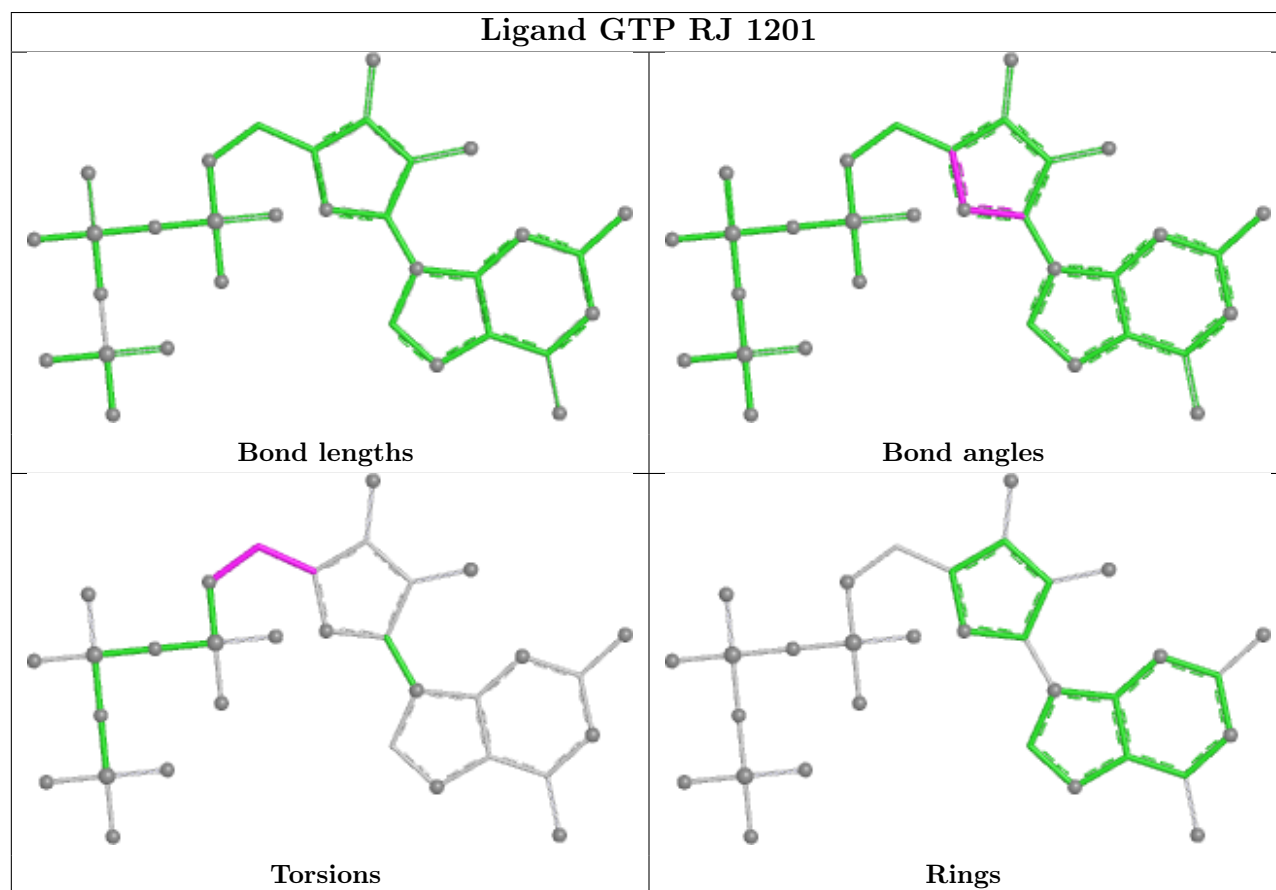
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
68	RJ	1201	GTP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

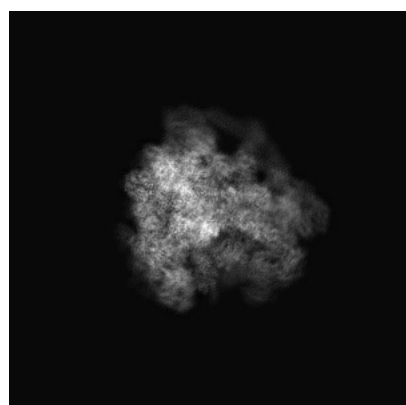
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0950. These allow visual inspection of the internal detail of the map and identification of artifacts.

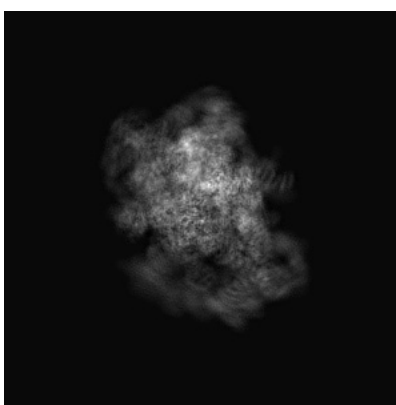
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

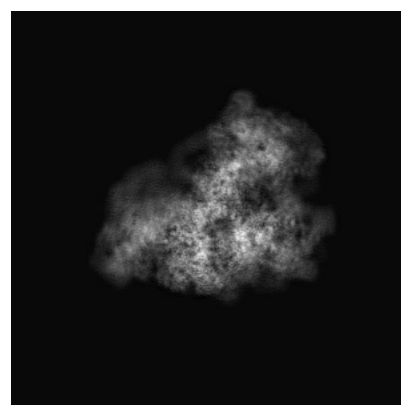
6.1.1 Primary map



X



Y

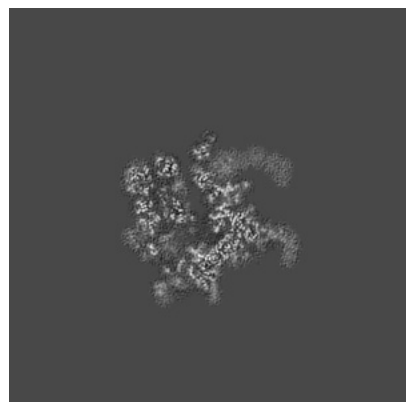


Z

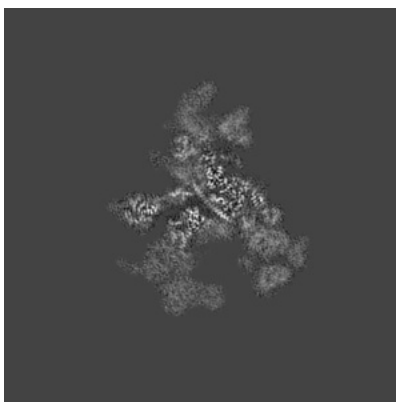
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

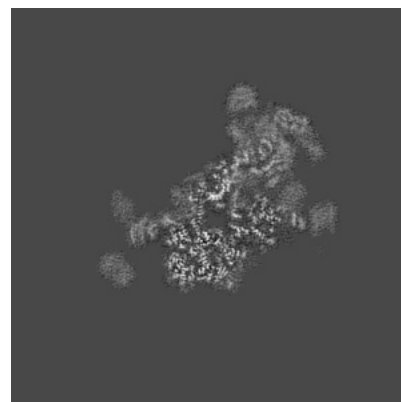
6.2.1 Primary map



X Index: 224



Y Index: 224

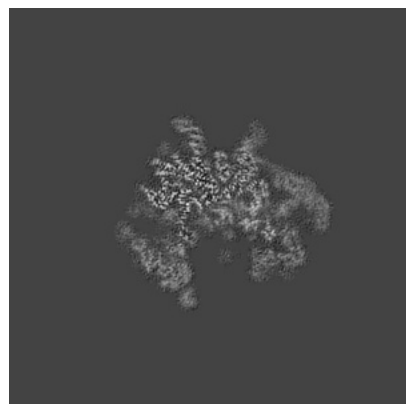


Z Index: 224

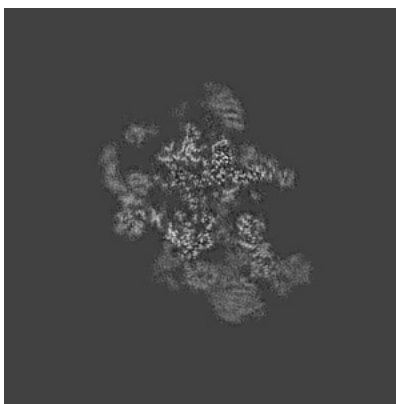
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

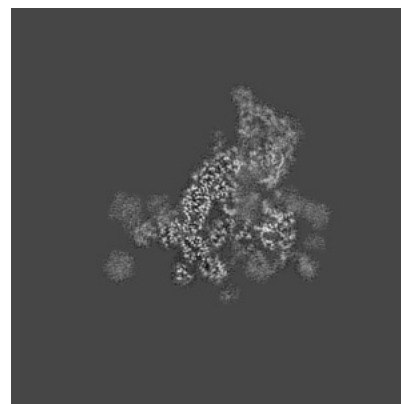
6.3.1 Primary map



X Index: 253



Y Index: 197

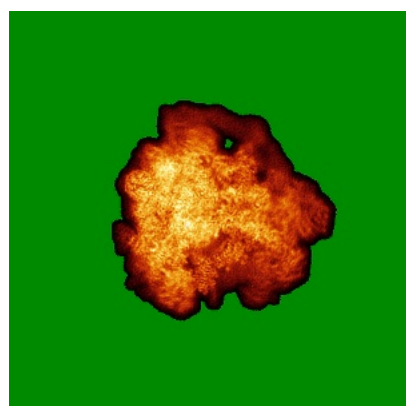


Z Index: 208

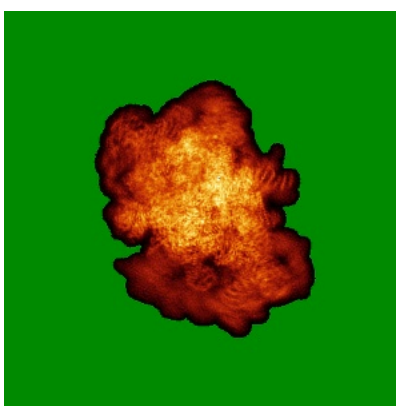
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

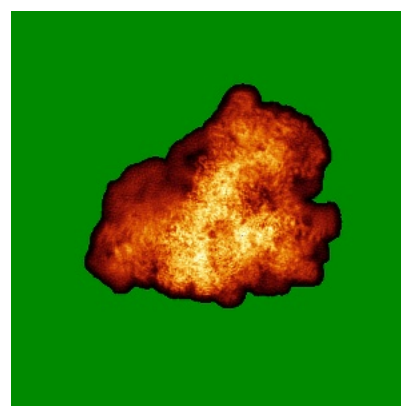
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

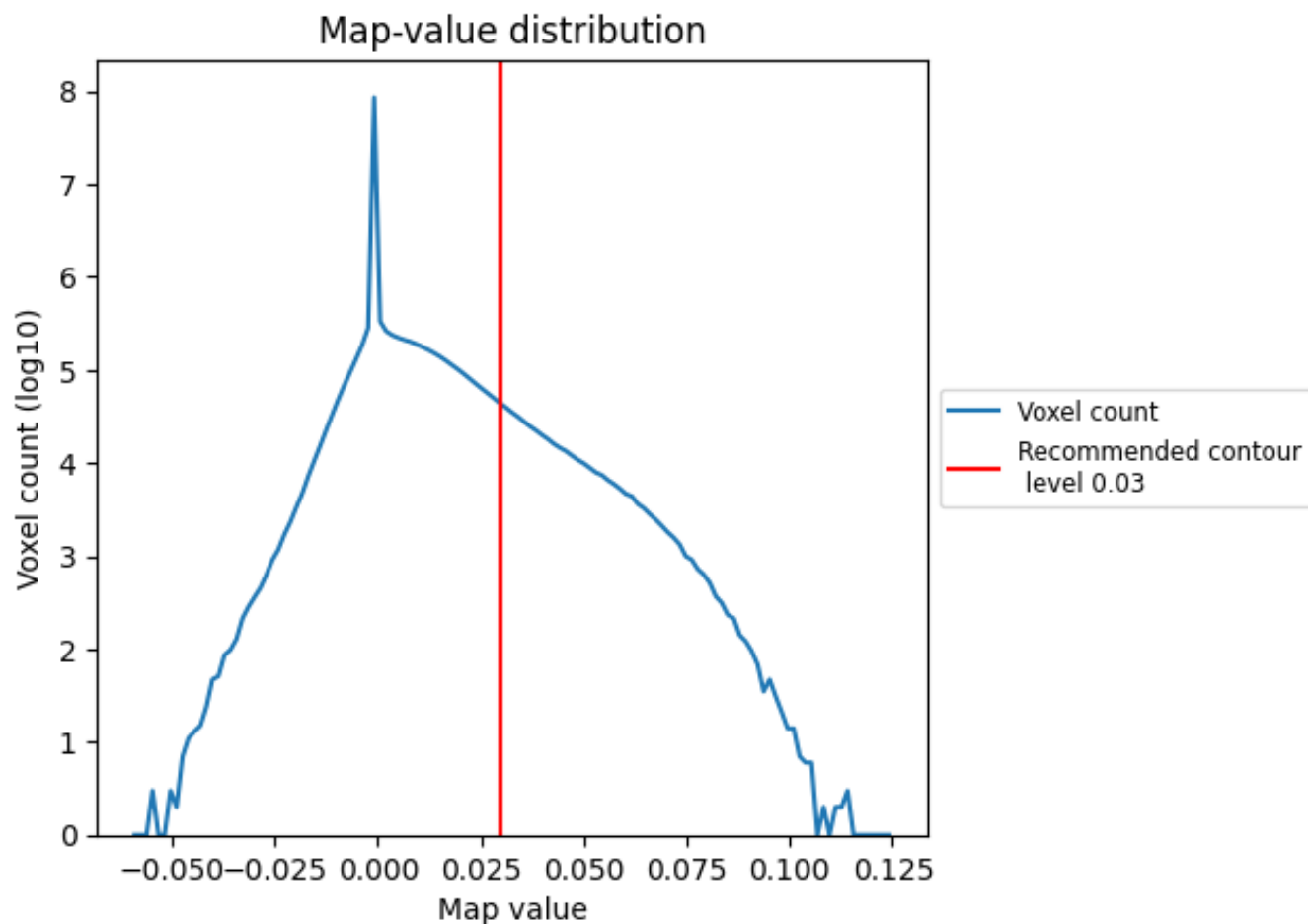
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

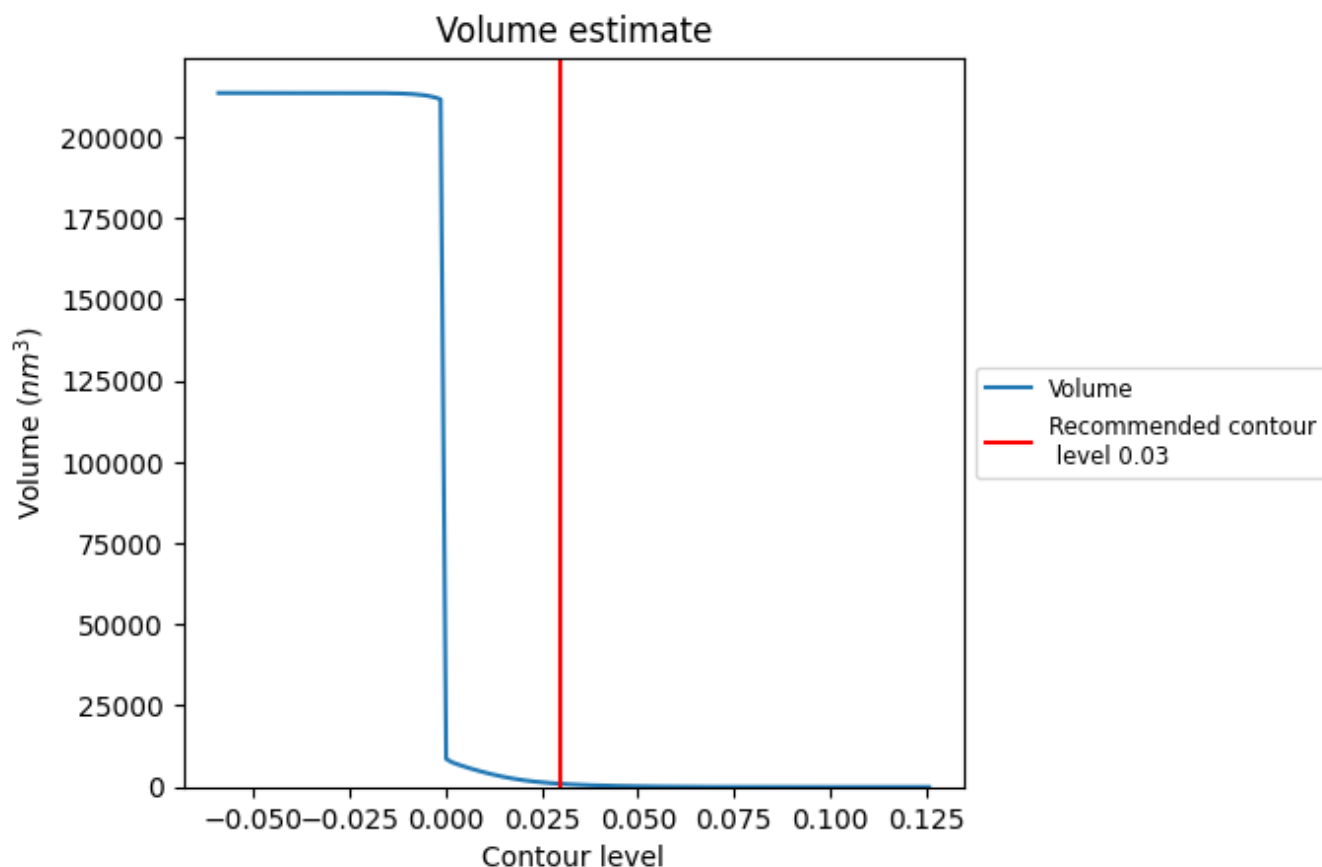
This section contains the results of statistical analysis of the map.

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.

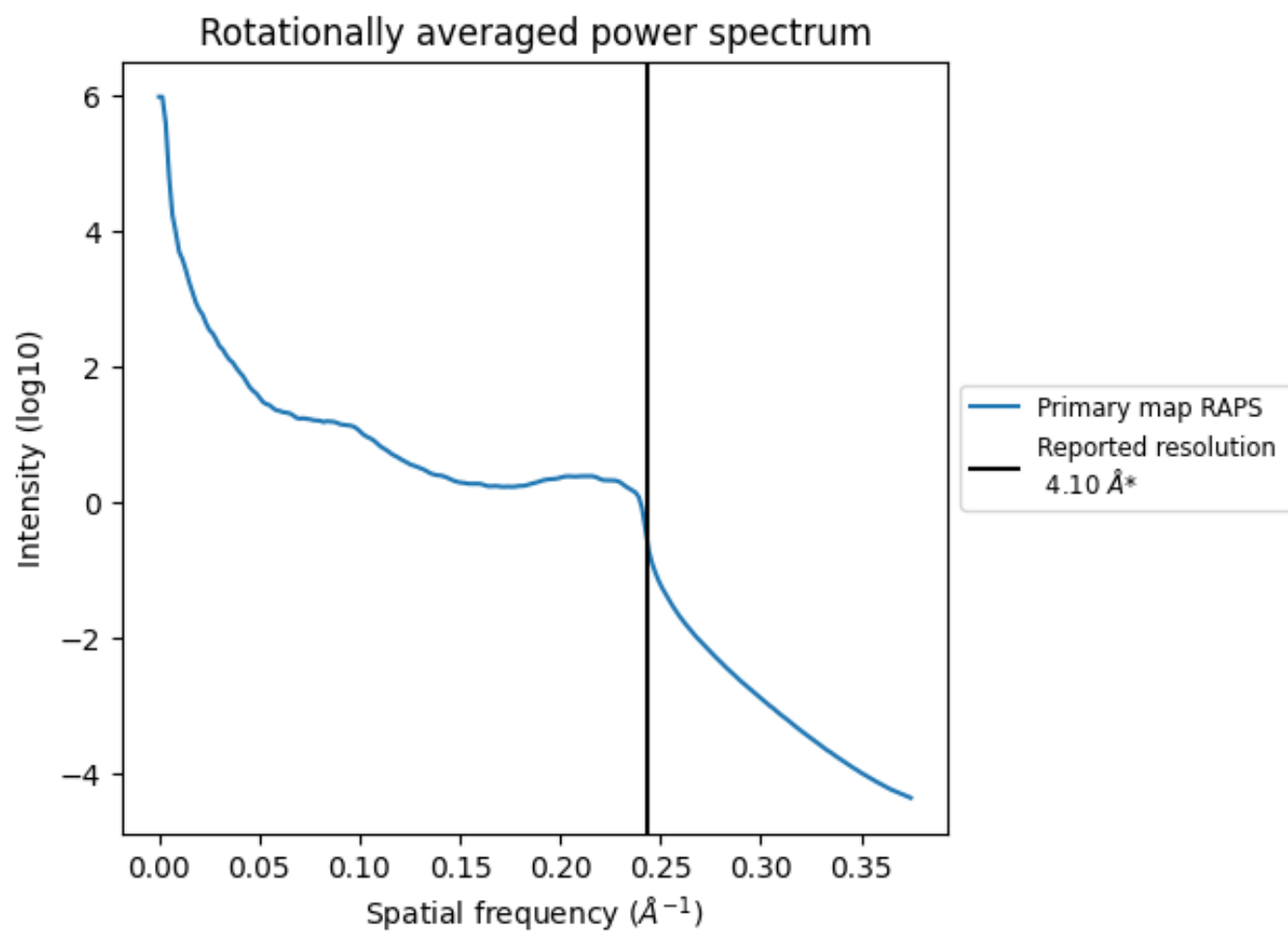
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 963 nm^3 ; this corresponds to an approximate mass of 870 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 ⓘ



*Reported resolution corresponds to spatial frequency of 0.244 Å⁻¹

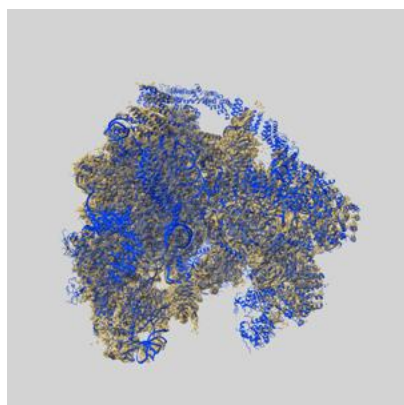
8 Fourier-Shell correlation

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

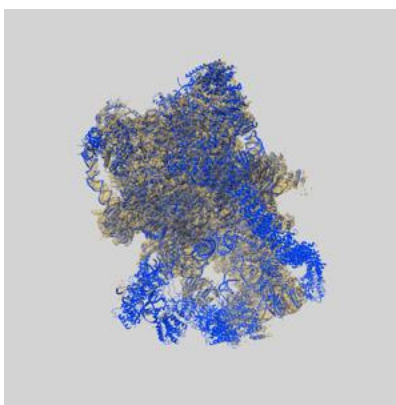
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-0950 and PDB model 6LQQ. Per-residue inclusion information can be found in section 3 on page 17.

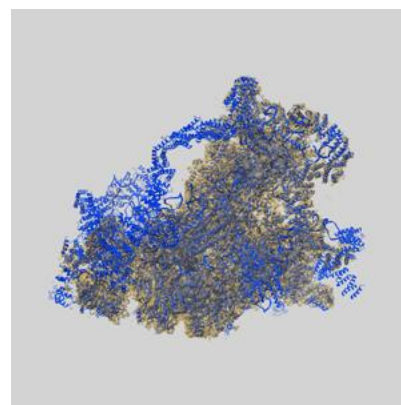
9.1 Map-model overlay [i](#)



X



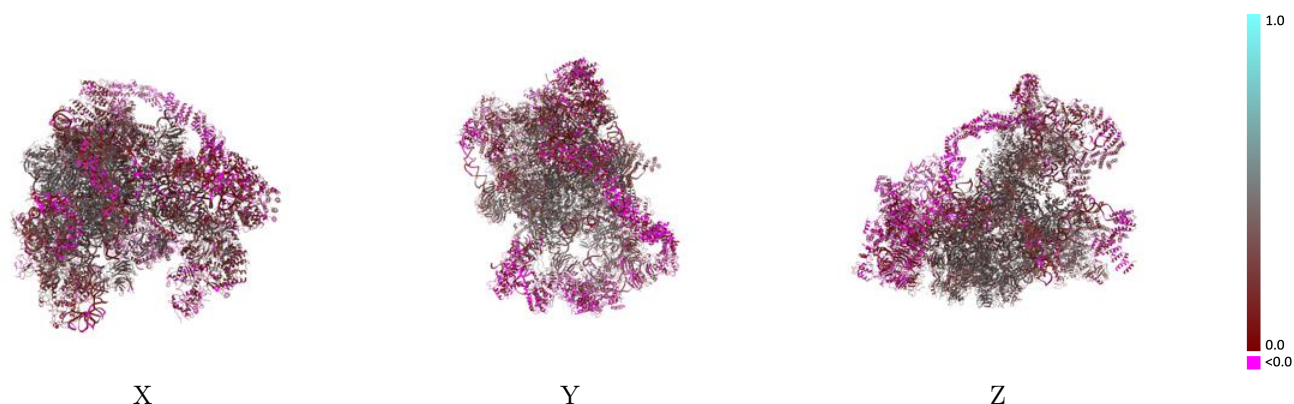
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.03 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](#)

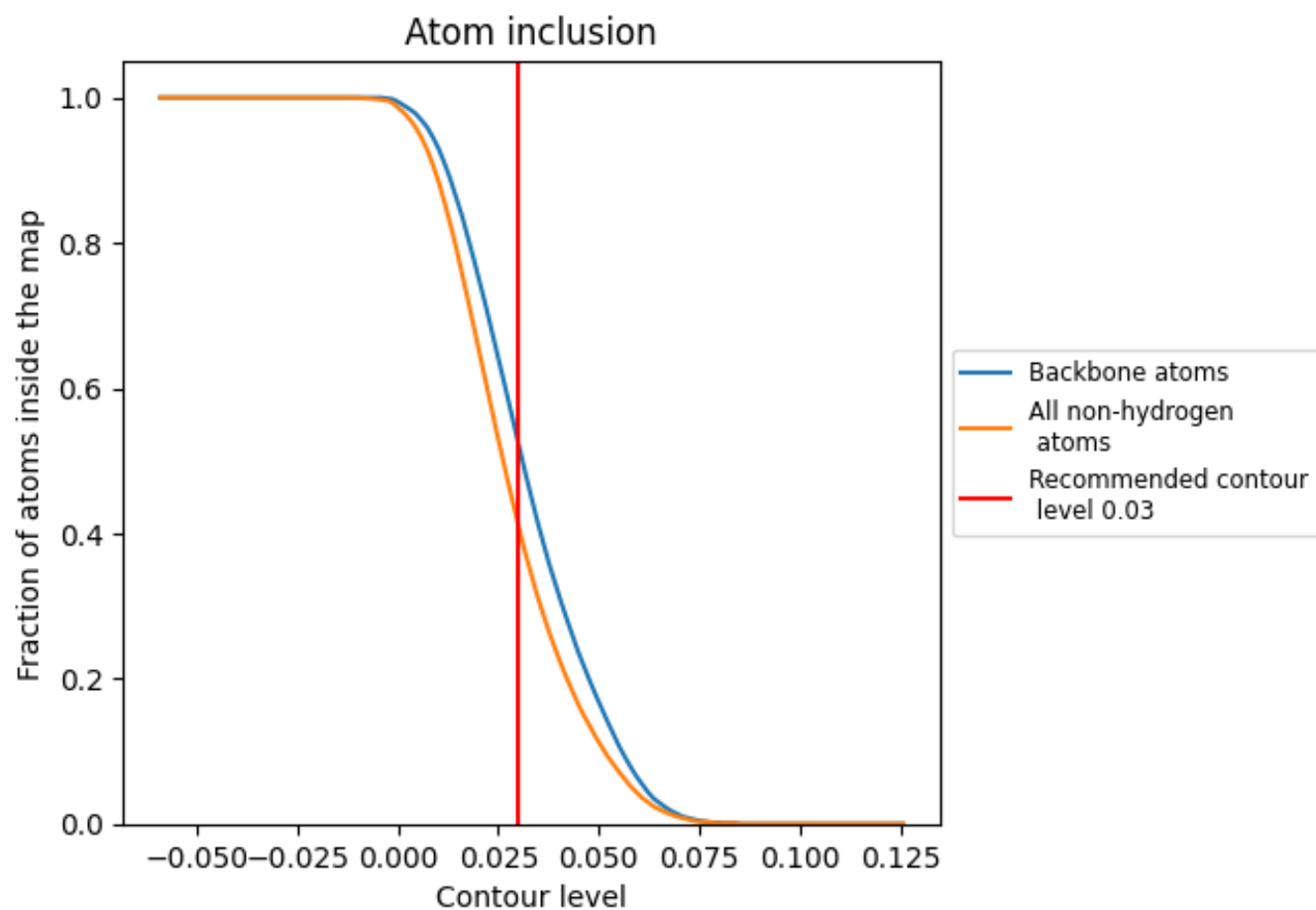


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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](#)

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 52% of all backbone atoms, 41% 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 (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.4110	 0.2550
3A	 0.6470	 0.2880
3B	 0.6400	 0.3850
3C	 0.1330	 0.1640
3D	 0.5950	 0.3190
3E	 0.4170	 0.2480
3F	 0.5760	 0.3330
3G	 0.5180	 0.2880
3H	 0.6120	 0.3680
5A	 0.2360	 0.2060
5B	 0.0290	 0.1290
5C	 0.5690	 0.3810
5D	 0.2920	 0.2640
5E	 0.5310	 0.3520
5F	 0.6520	 0.3990
5G	 0.5730	 0.3840
5H	 0.6500	 0.3910
5I	 0.6900	 0.4090
5J	 0.3860	 0.2600
5K	 0.6320	 0.4030
A4	 0.2650	 0.1750
A5	 0.4090	 0.2640
A8	 0.0170	 0.0350
A9	 0.0500	 0.0530
AE	 0.2040	 0.1300
AF	 0.1830	 0.1930
AG	 0.2310	 0.1370
B1	 0.6490	 0.4020
B2	 0.5870	 0.3260
B3	 0.4860	 0.2820
B6	 0.5520	 0.2760
B8	 0.5770	 0.3310
BE	 0.6500	 0.3950
RA	 0.4000	 0.2630
RB	 0.4710	 0.2910



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Chain	Atom inclusion	Q-score
RC	 0.3940	 0.3180
RD	 0.0020	 0.0240
RE	 0.3440	 0.2480
RF	 0.2870	 0.2400
RG	 0.0150	 0.1160
RH	 0.0410	 0.1740
RJ	 0.5980	 0.3620
RK	 0.6270	 0.3690
RL	 0.4210	 0.2740
RM	 0.1620	 0.1440
RN	 0.0760	 0.1140
RO	 0.0000	 0.0830
RP	 0.2500	 0.1580
RQ	 0.4550	 0.3130
RS	 0.0020	 0.0210
RT	 0.2210	 0.3060
RV	 0.3570	 0.3440
RY	 0.3040	 0.2060
SA	 0.5390	 0.2470
SC	 0.5380	 0.3630
SF	 0.4940	 0.2910
SG	 0.6060	 0.3600
SH	 0.3660	 0.2430
SI	 0.4620	 0.2910
SJ	 0.3960	 0.2340
SK	 0.6400	 0.3810
SM	 0.2530	 0.1500
SO	 0.5920	 0.3670
SP	 0.5510	 0.3690
SR	 0.6340	 0.3890
SX	 0.5970	 0.3850
SY	 0.6040	 0.4080
SZ	 0.5550	 0.3210
Sc	 0.5900	 0.3860
Sd	 0.5950	 0.3980
X1	 0.0830	 0.1700