



## wwPDB EM Validation Summary Report ⓘ

Nov 13, 2022 – 02:04 PM EST

PDB ID : 6VYZ  
EMDB ID : EMD-21477  
Title : Escherichia coli transcription-translation complex C6 (TTC-C6) containing mRNA with a 21 nt long spacer, NusA, and fMet-tRNAs at E-site and P-site  
Authors : Molodtsov, V.; Wang, C.; Su, M.; Ebright, R.H.  
Deposited on : 2020-02-27  
Resolution : 9.90 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

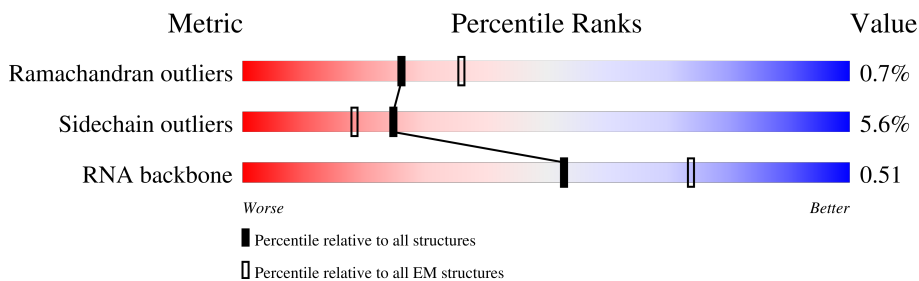
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

# 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 9.90 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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	0	103	
2	1	110	
3	2	100	
4	3	104	
5	4	94	
6	5	36	
7	6	36	
8	7	38	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
9	A	76	25% 58% 38%
9	B	76	18% 47% 49%
10	AA	1342	34% 86% 9%
11	AB	181	33% 53% 46%
12	AC	329	35% 65% 5% 30%
12	AD	329	34% 69% 31%
13	AE	1407	42% 89% 5% 5%
14	AF	91	65% 90% 9%
15	C	75	11% 85% 12%
16	D	1542	14% 79% 20%
17	E	87	46% 93% 6%
18	F	71	76% 94%
19	G	241	18% 90% 7%
20	H	557	31% 41% 54%
21	I	208	33% 97%
22	J	206	42% 97%
23	K	167	40% 89% 5% 7%
24	L	135	11% 72% 23%
25	M	179	27% 80% 16%
26	N	130	39% 97%
27	O	130	36% 93% 5%
28	P	103	35% 89% 7%
29	Q	129	19% 87% 9%
30	R	124	30% 92% 6%
31	S	101	42% 95%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
32	T	89	17% 85% 13%
33	U	82	54% 93% 7%
34	V	84	43% 93% 5%
35	W	92	17% 86% 10%
36	X	118	27% 88% 10%
37	Y	3	100% 33% 67%
38	a	2904	21% 81% 18%
39	b	85	48% 88% 11%
40	c	78	51% 94% 5%
41	d	120	38% 86% 14%
42	e	62	32% 98%
43	f	59	64% 95%
44	g	70	10% 86% 9% 6%
45	h	273	43% 93% 7%
46	i	57	46% 88% 11%
47	j	209	48% 97%
48	k	55	25% 89% 5% 5%
49	l	201	38% 93% 7%
50	m	46	61% 93% 7%
51	n	179	16% 89% 10%
52	o	65	86% 91% 8%
53	p	177	17% 97%
54	q	38	37% 95% 5%
55	r	149	82% 93% 7%
56	s	142	57% 96%

Continued on next page...

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
57	t	123	<p>43% 95% 5%</p>
58	u	144	<p>42% 96%</p>
59	v	136	<p>38% 96%</p>
60	w	127	<p>51% 87% 6% 6%</p>
61	x	117	<p>27% 94% 5%</p>
62	y	115	<p>57% 95%</p>
63	z	118	<p>44% 97%</p>

## 2 Entry composition [i](#)

There are 65 unique types of molecules in this entry. The entry contains 297741 atoms, of which 124644 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 protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
1	0	103	1655	516	839	153	145	2	0	0

- Molecule 2 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	1	110	1779	532	922	166	156	3	0	0

- Molecule 3 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
3	2	94	1557	470	811	140	134	2	0	0

- Molecule 4 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
4	3	103	1632	498	844	148	142	0	0

- Molecule 5 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
5	4	94	1533	479	780	137	134	3	0	0

- Molecule 6 is a DNA chain called NT DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
6	5	23	731	225	259	87	137	23	0	0

- Molecule 7 is a DNA chain called T DNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
7	6	27	847	259	305	89	167	27	0	0

- Molecule 8 is a RNA chain called mRNA with 21 nt long spacer.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
8	7	16	515	154	168	62	115	16	0	0

- Molecule 9 is a RNA chain called E-site and P-site tRNA (fMet).

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
9	A	76	2446	723	826	295	527	75	0	0
9	B	76	2433	723	813	295	527	75	0	0

- Molecule 10 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
10	AA	1223	19199	6029	9589	1672	1867	42	0	0

- Molecule 11 is a protein called Transcription termination/antitermination protein NusG.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
11	AB	98	1573	505	783	139	140	6	0	0

- Molecule 12 is a protein called DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
12	AC	230	3599	1112	1813	317	351	6	0	0
12	AD	228	3556	1100	1789	312	349	6	0	0

- Molecule 13 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
13	AE	1335	20999	6526	10611	1854	1958	50	0	0

- Molecule 14 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
14	AF	83	1318	399	663	123	132	1	0	0

- Molecule 15 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
15	C	66	1103	344	559	102	97	1	0	0

- Molecule 16 is a RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
16	D	1522	49076	14565	16420	5993	10576	1522	0	0

- Molecule 17 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
17	E	86	1388	414	719	138	114	3	0	0

- Molecule 18 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
18	F	70	1218	366	629	125	97	1	0	0

- Molecule 19 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
19	G	225	3545	1113	1785	316	323	8	0	0

- Molecule 20 is a protein called 30S ribosomal protein S1.



Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
20	H	259	3184	1073	1454	305	349	3	0	0

- Molecule 21 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
21	I	208	3346	1036	1710	307	290	3	0	0

- Molecule 22 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
22	J	205	3350	1026	1707	315	298	4	0	0

- Molecule 23 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
23	K	156	2348	717	1196	217	212	6	0	0

- Molecule 24 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
24	L	104	1694	536	846	153	152	7	0	0

- Molecule 25 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
25	M	151	2416	735	1235	227	215	4	0	0

- Molecule 26 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
26	N	129	2010	616	1031	173	184	6	0	0

- Molecule 27 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
27	O	127	2092	634	1070	206	179	3	0	0

- Molecule 28 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
28	P	99	1621	495	831	151	143	1	0	0

- Molecule 29 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
29	Q	117	1764	540	887	174	160	3	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
30	R	121	1940	580	1001	194	161	4	0	0

- Molecule 31 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
31	S	100	1649	499	844	164	139	3	0	0

- Molecule 32 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
32	T	88	1448	439	734	144	130	1	0	0

- Molecule 33 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
33	U	82	1315	406	666	128	114	1	0	0

- Molecule 34 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
34	V	80	1339	411	691	121	113	3	0	0

- Molecule 35 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
35	W	83	1351	424	688	126	111	2	0	0

- Molecule 36 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
36	X	116	1864	558	964	181	158	3	0	0

- Molecule 37 is a RNA chain called mRNA in the ribosomal RNA entrance pore.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
37	Y	3	90	27	30	6	24	3	0	0

- Molecule 38 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
38	a	2880	92917	27587	31076	11398	19976	2880	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	887	A	U	conflict	GB 937521852

- Molecule 39 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
39	b	76	1181	360	599	117	104	1	0	0

- Molecule 40 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
40	c	77	1277	388	652	129	106	2	0	0

- Molecule 41 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
41	d	120	3870	1144	1301	468	837	120	0	0

- Molecule 42 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
42	e	62	1032	308	531	98	94	1	0	0

- Molecule 43 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
43	f	58	936	281	488	87	78	2	0	0

- Molecule 44 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
44	g	66	1042	323	520	99	94	6	0	0

- Molecule 45 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
45	h	271	4236	1288	2154	423	364	7	0	0

- Molecule 46 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
46	i	56	903	269	459	94	80	1	0	0

- Molecule 47 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
47	j	209	Total	C	H	N	O	S	0	0
			3182	979	1617	288	294	4		

- Molecule 48 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					AltConf	Trace	
48	k	52	Total	C	H	N	O	S	0	0
			890	275	464	78	73			

- Molecule 49 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace	
49	l	201	Total	C	H	N	O	S	0	0
			3171	974	1619	283	290	5		

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

Mol	Chain	Residues	Atoms					AltConf	Trace	
50	m	46	Total	C	H	N	O	S	0	0
			795	228	418	90	57	2		

- Molecule 51 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace	
51	n	177	Total	C	H	N	O	S	0	0
			2853	899	1443	249	256	6		

- Molecule 52 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace	
52	o	64	Total	C	H	N	O	S	0	0
			1076	323	572	105	74	2		

- Molecule 53 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace	
53	p	175	Total	C	H	N	O	S	0	0
			2671	826	1358	241	244	2		

- Molecule 54 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
54	q	38	645	185	343	65	48	4	0	0

- Molecule 55 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
55	r	149	2259	699	1148	197	214	1	0	0

- Molecule 56 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
56	s	142	2291	714	1162	212	199	4	0	0

- Molecule 57 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
57	t	123	1969	593	1023	181	166	6	0	0

- Molecule 58 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
58	u	144	2182	654	1129	207	190	2	0	0

- Molecule 59 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
59	v	136	2231	686	1157	205	177	6	0	0

- Molecule 60 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
60	w	119	1945	588	994	195	163	5	0	0

- Molecule 61 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
61	x	116	1815	552	923	178	162	0	0

- Molecule 62 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
62	y	114	1879	574	962	179	163	1	0	0

- Molecule 63 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
63	z	117	1967	604	1020	192	151	0	0

- Molecule 64 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
64	AE	1	1	1	0

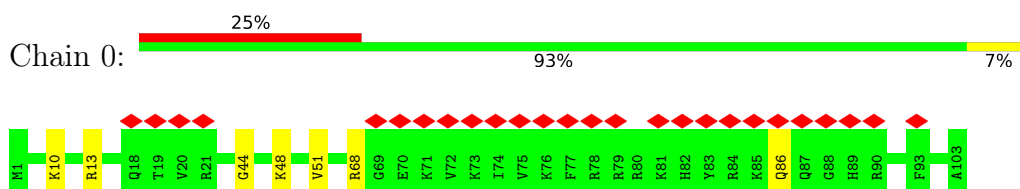
- Molecule 65 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
			Total	Zn	
65	AE	2	2	2	0

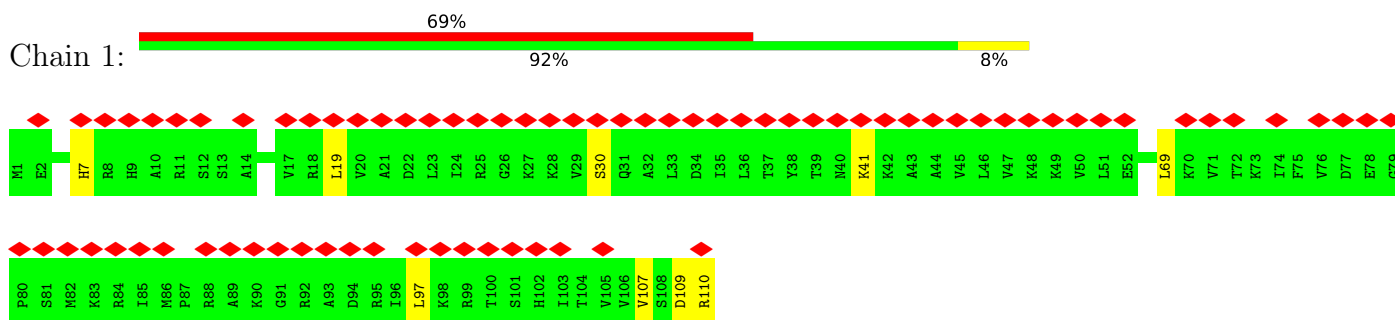
### 3 Residue-property plots [i](#)

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.

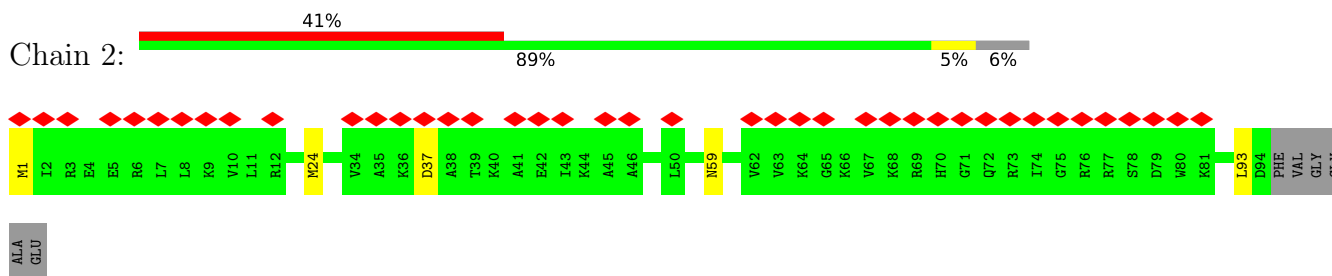
- Molecule 1: 50S ribosomal protein L21



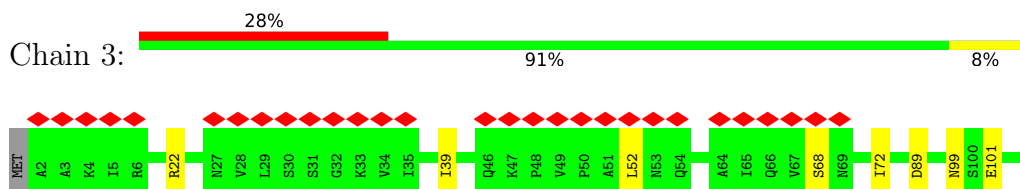
- Molecule 2: 50S ribosomal protein L22



- Molecule 3: 50S ribosomal protein L23



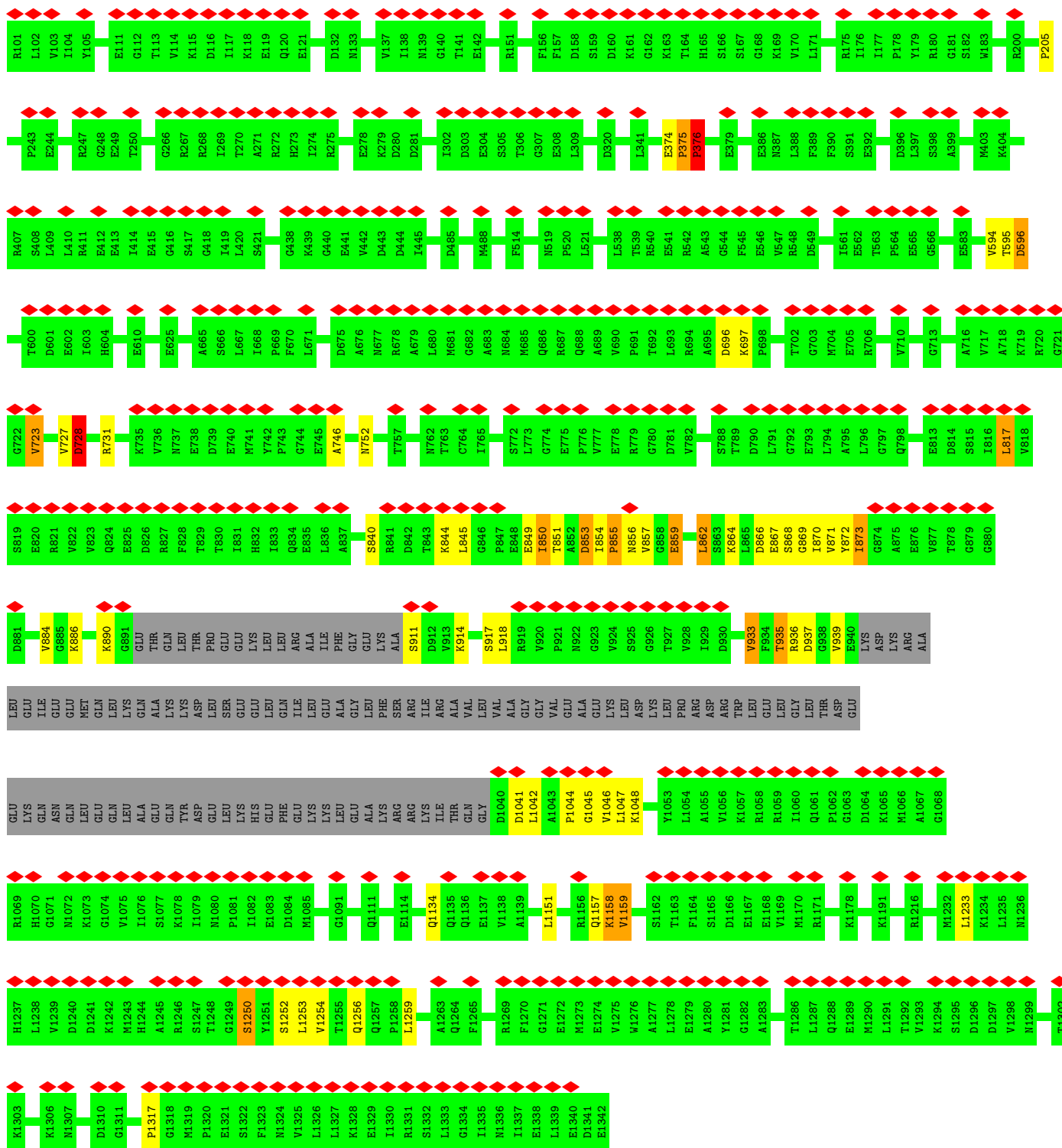
- Molecule 4: 50S ribosomal protein L24



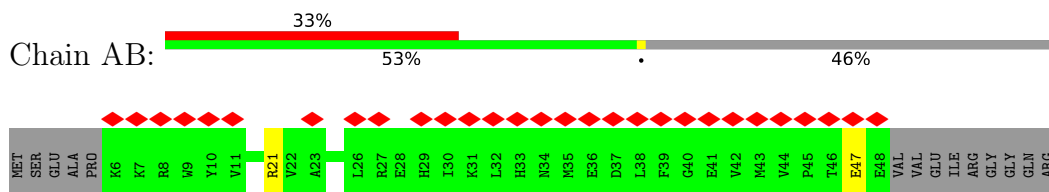
- Molecule 5: 50S ribosomal protein L25

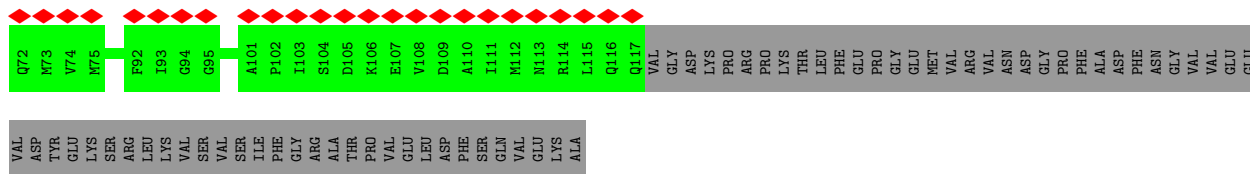




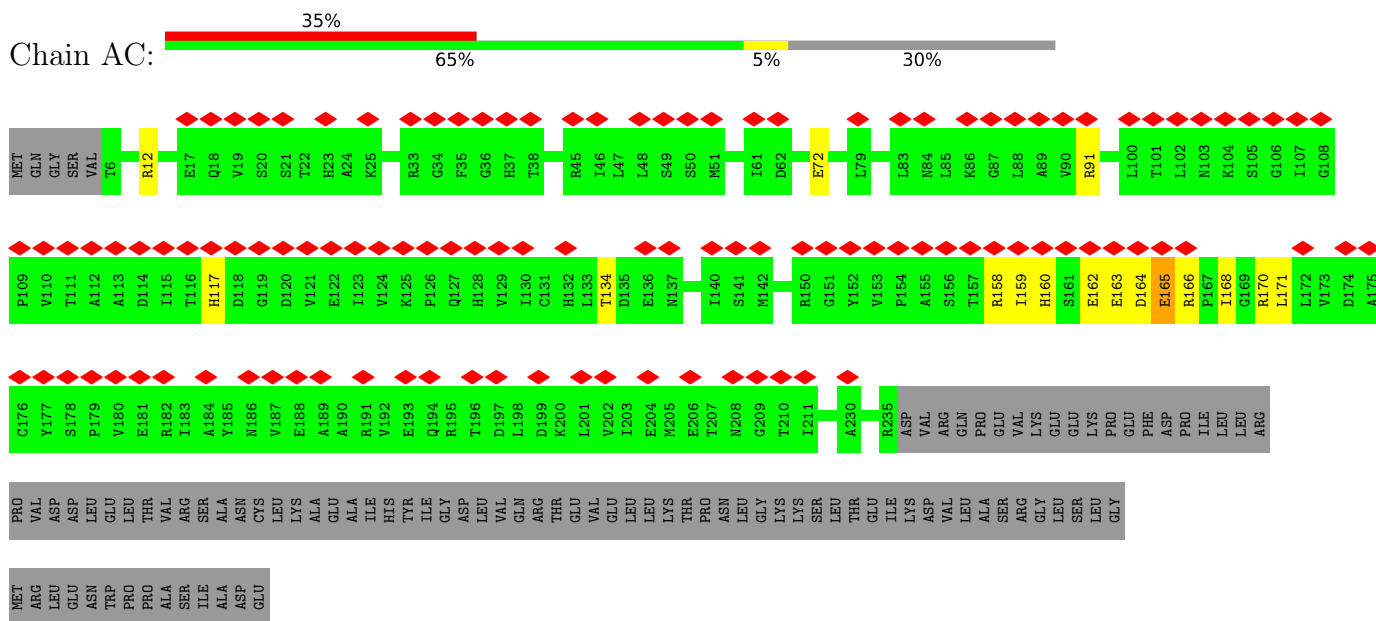


• Molecule 11: Transcription termination/antitermination protein NusG

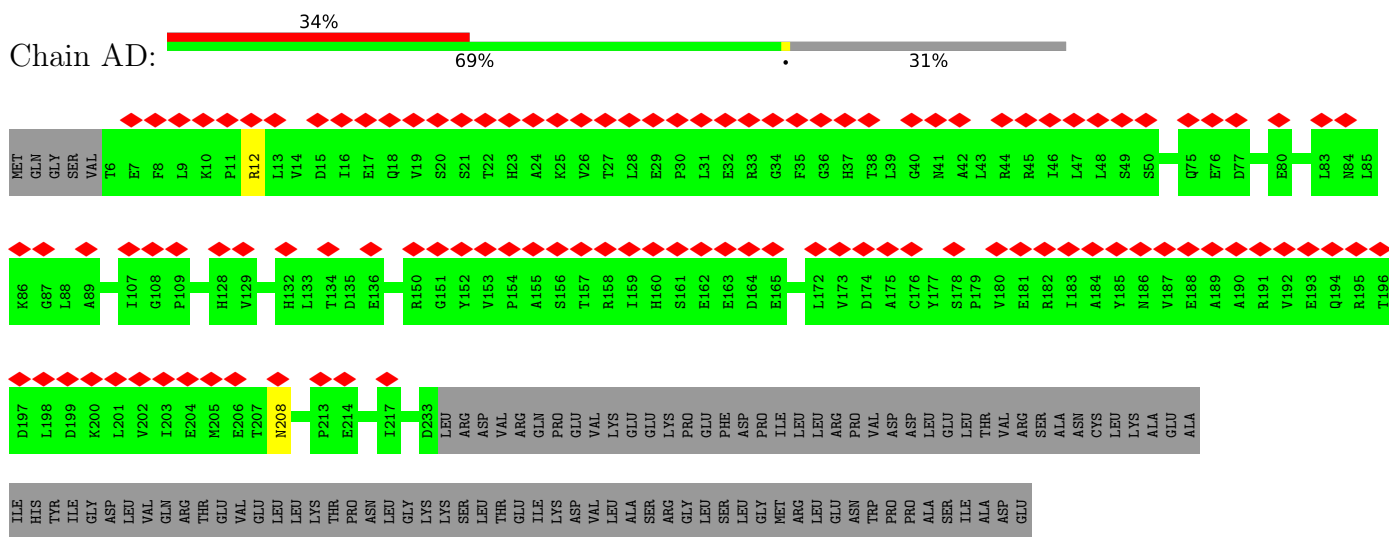




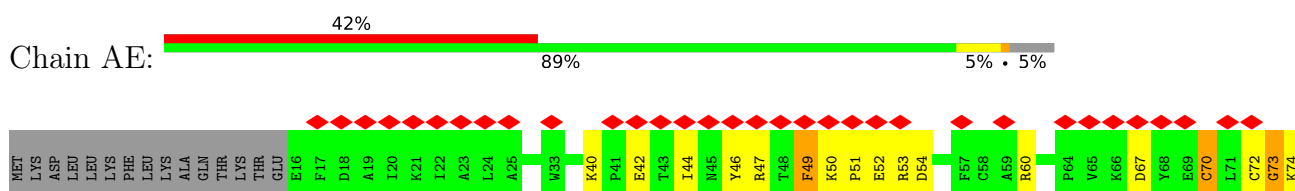
● Molecule 12: DNA-directed RNA polymerase subunit alpha



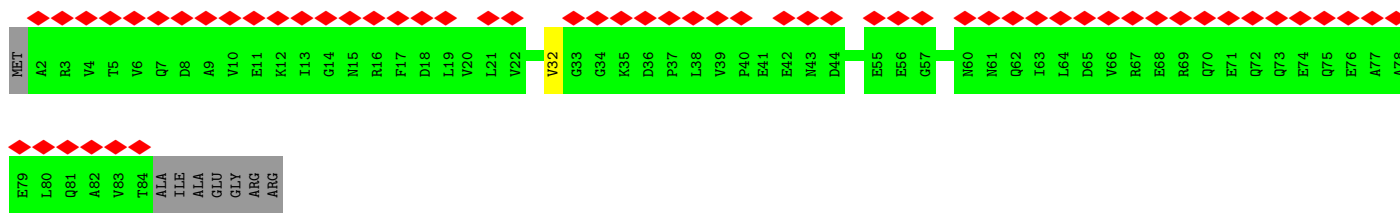
● Molecule 12: DNA-directed RNA polymerase subunit alpha



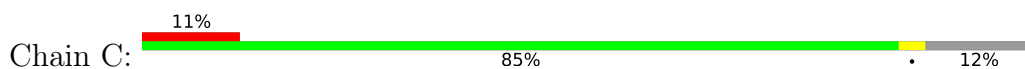
● Molecule 13: DNA-directed RNA polymerase subunit beta'



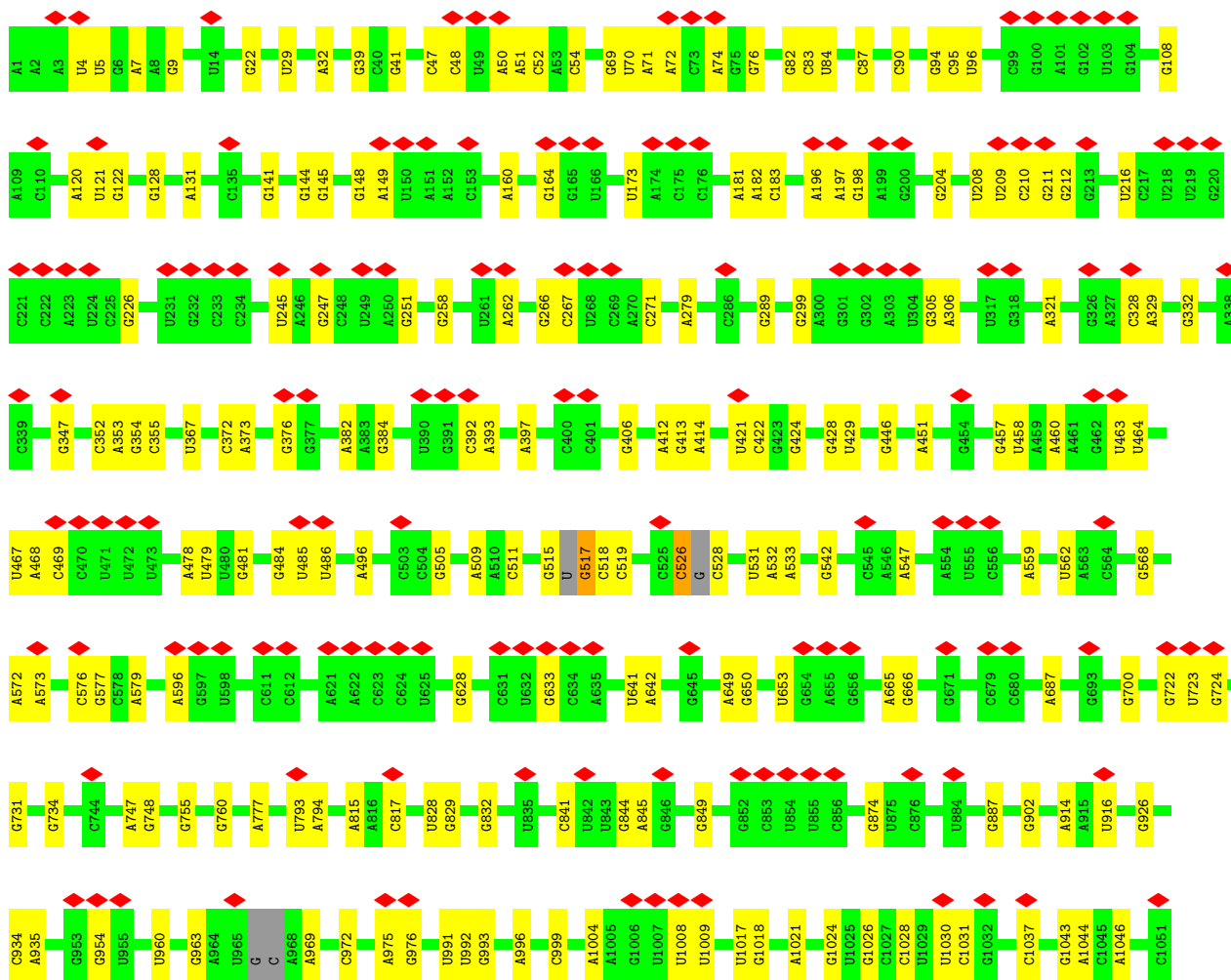
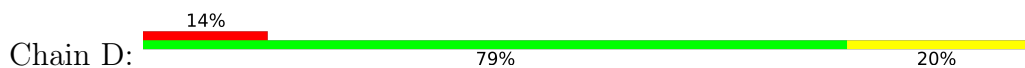


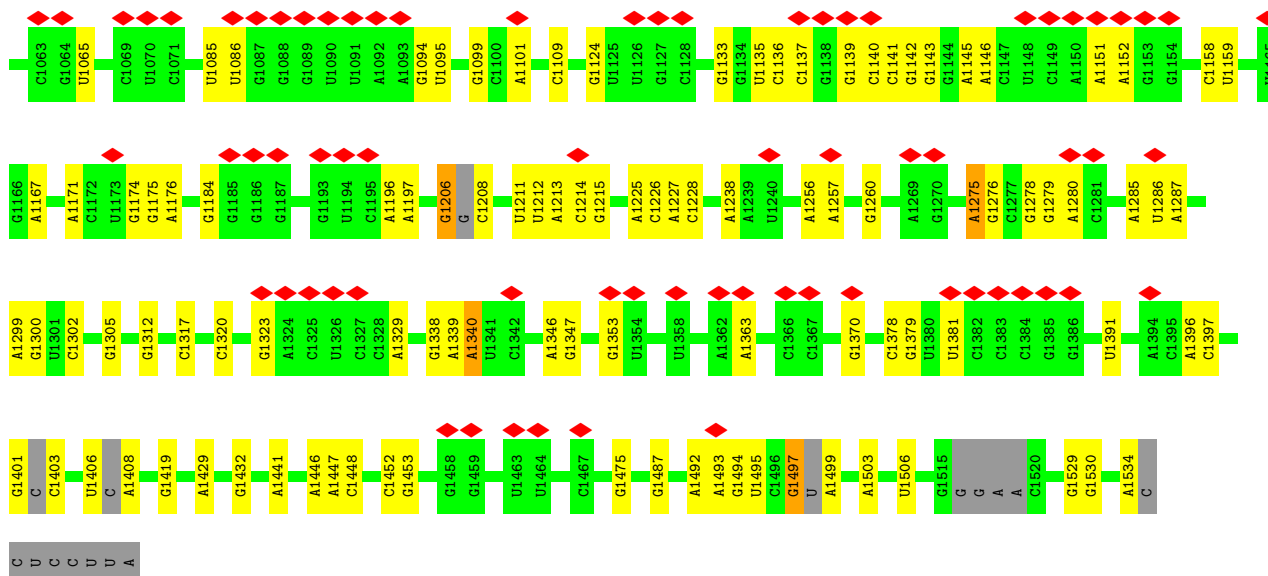


• Molecule 15: 30S ribosomal protein S18

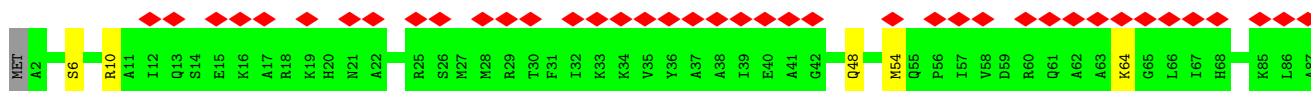
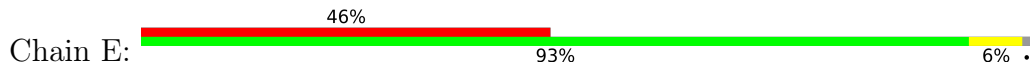


• Molecule 16: 16S rRNA

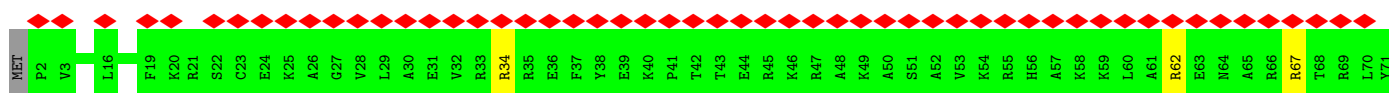
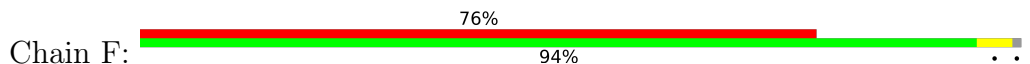




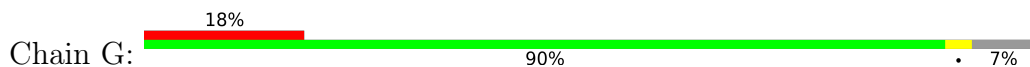
• Molecule 17: 30S ribosomal protein S20



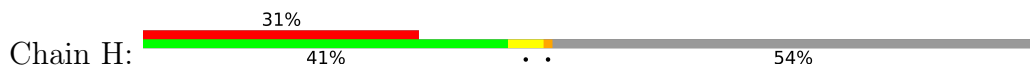
• Molecule 18: 30S ribosomal protein S21



• Molecule 19: 30S ribosomal protein S2



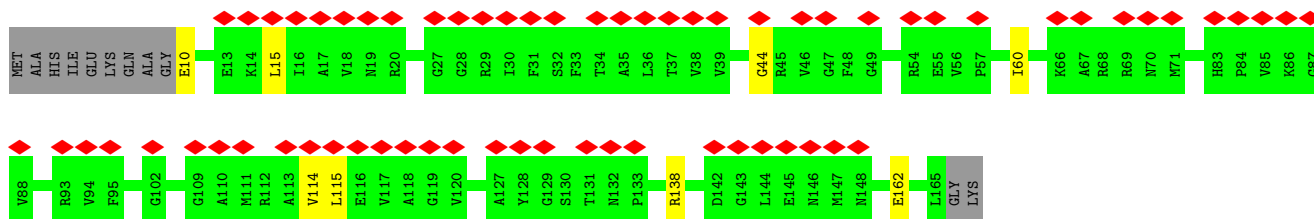
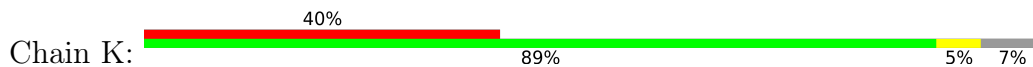
• Molecule 20: 30S ribosomal protein S1



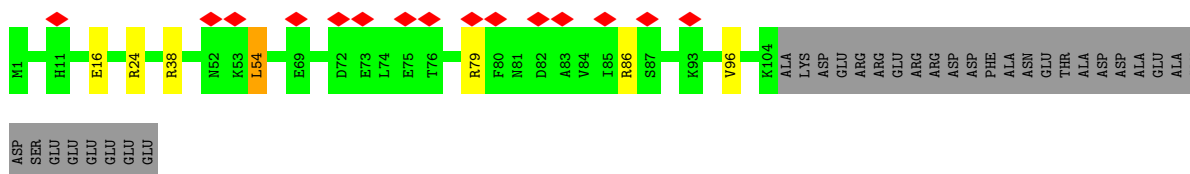
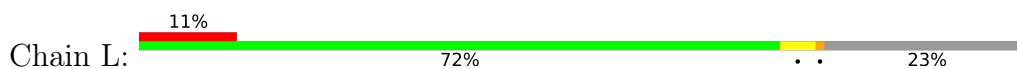


◆  
K206

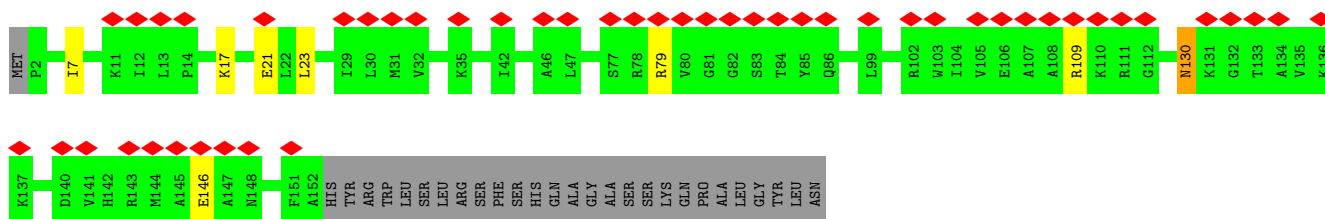
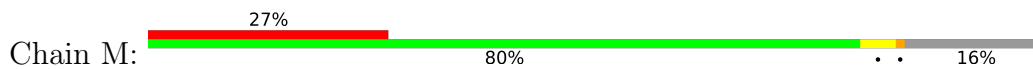
• Molecule 23: 30S ribosomal protein S5



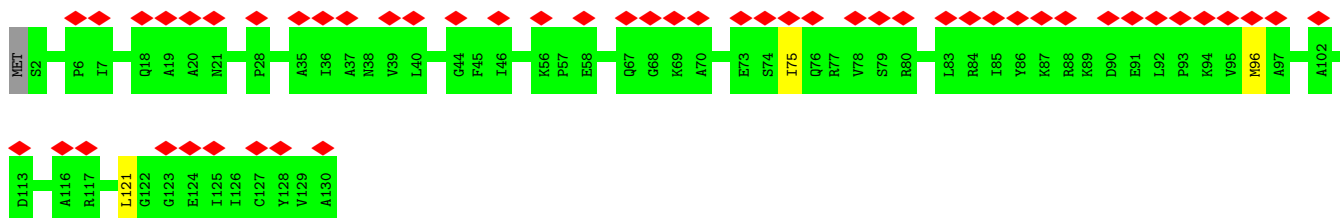
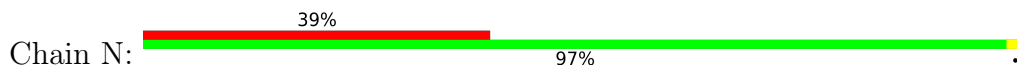
• Molecule 24: 30S ribosomal protein S6



• Molecule 25: 30S ribosomal protein S7



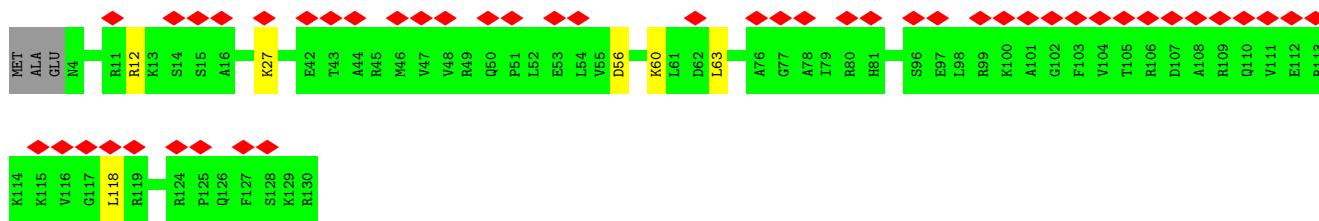
• Molecule 26: 30S ribosomal protein S8



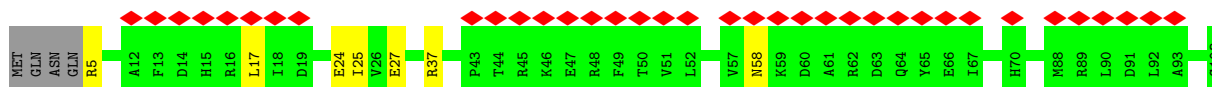
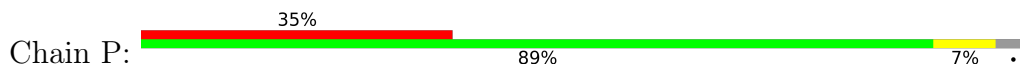
• Molecule 27: 30S ribosomal protein S9



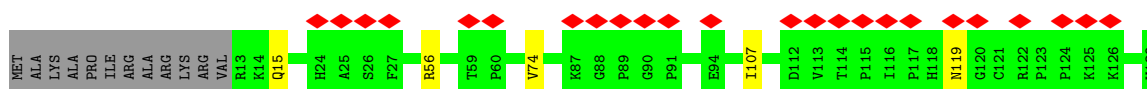
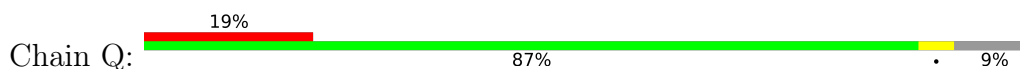




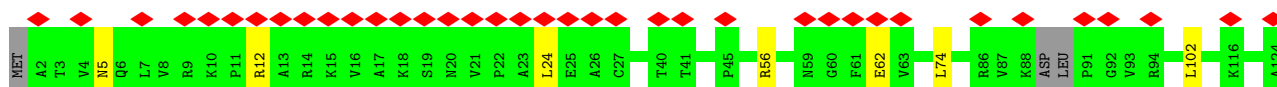
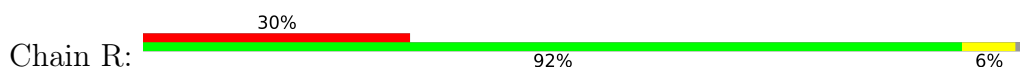
- Molecule 28: 30S ribosomal protein S10



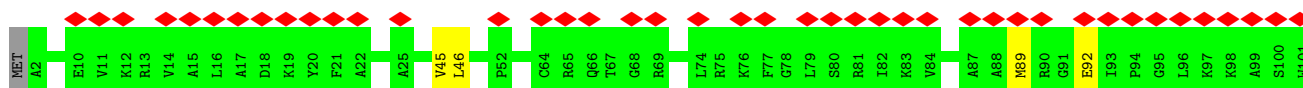
- Molecule 29: 30S ribosomal protein S11



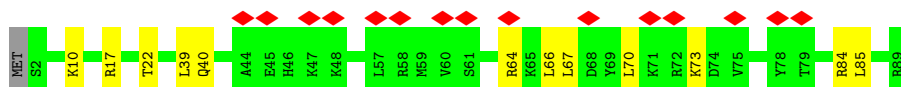
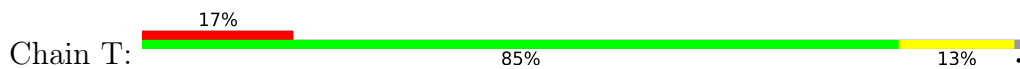
- Molecule 30: 30S ribosomal protein S12



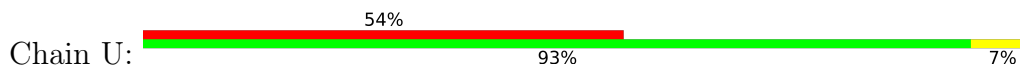
- Molecule 31: 30S ribosomal protein S14



- Molecule 32: 30S ribosomal protein S15

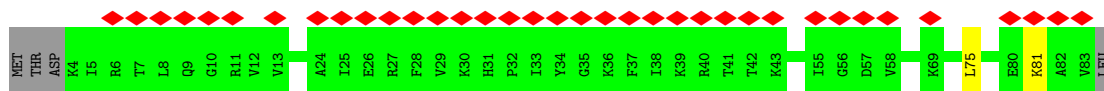
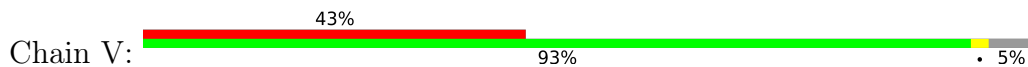


- Molecule 33: 30S ribosomal protein S16

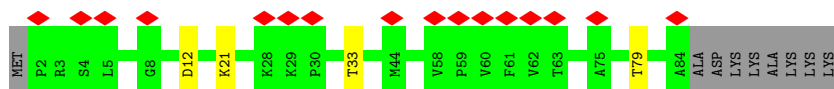
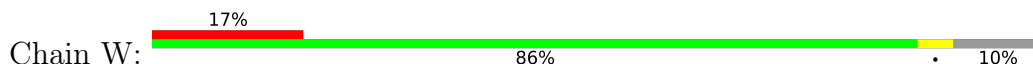




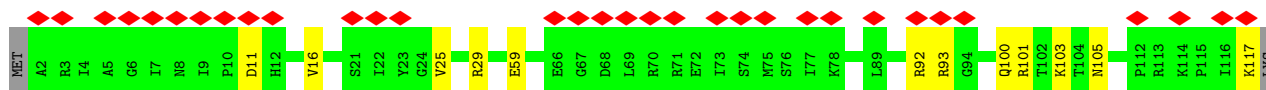
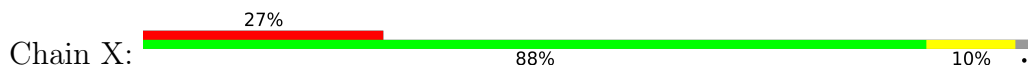
• Molecule 34: 30S ribosomal protein S17



• Molecule 35: 30S ribosomal protein S19



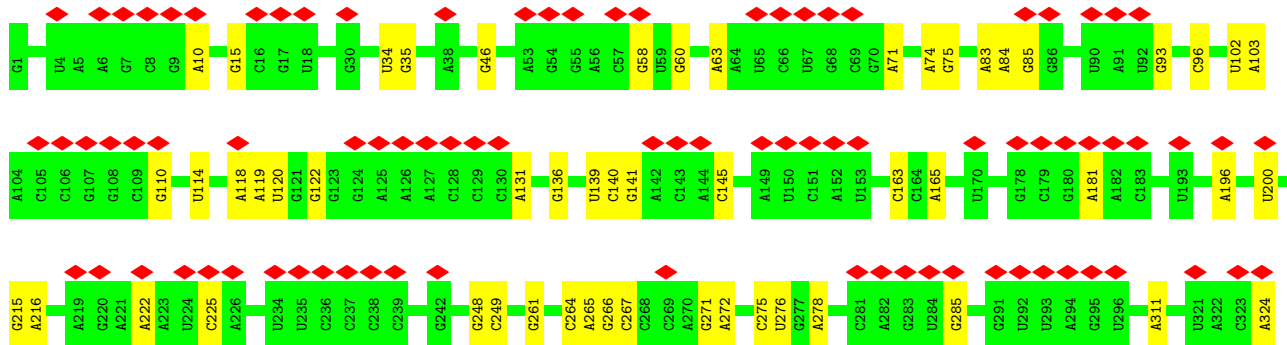
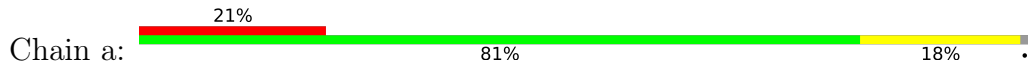
• Molecule 36: 30S ribosomal protein S13

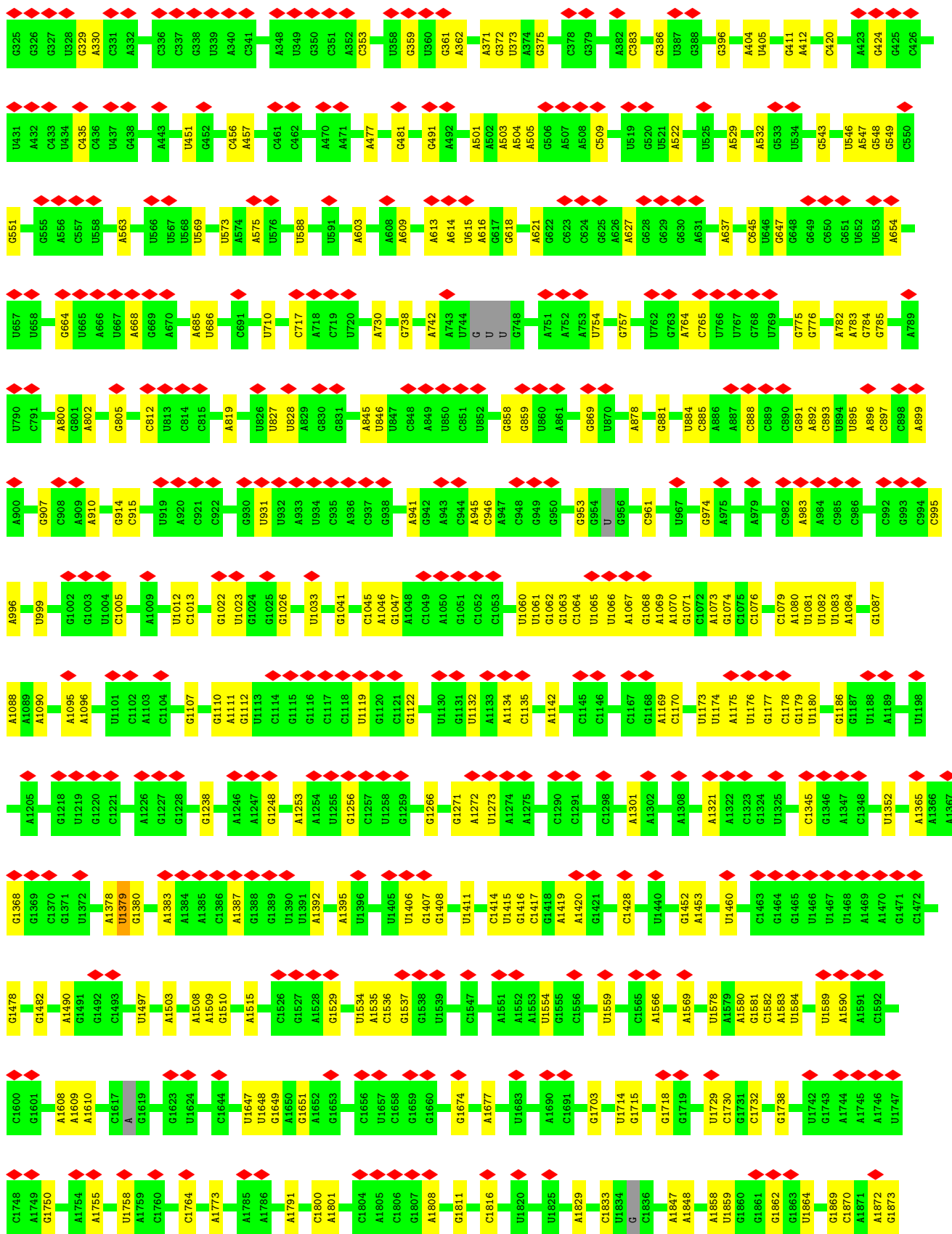


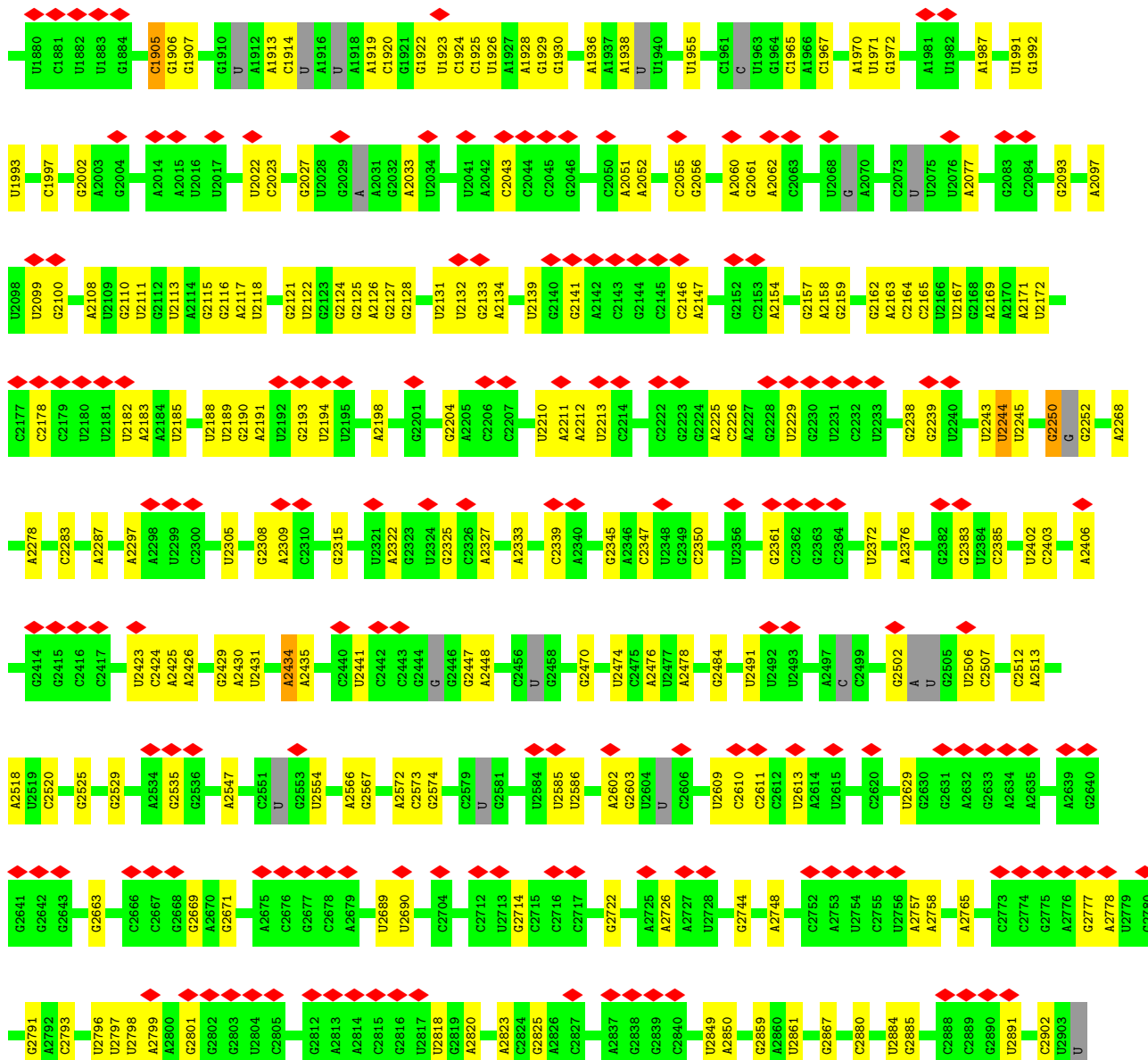
• Molecule 37: mRNA in the ribosomal RNA entrance pore



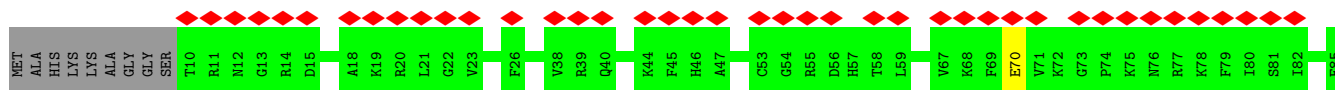
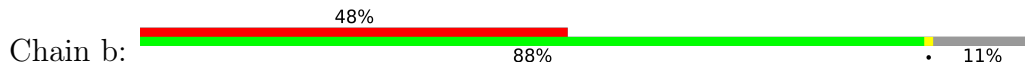
• Molecule 38: 23S rRNA



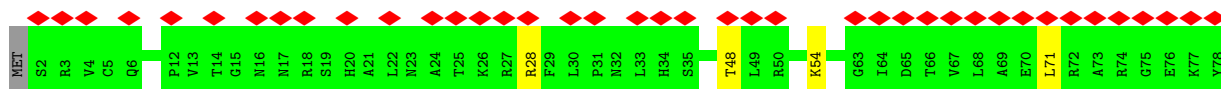
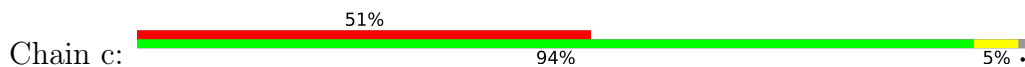




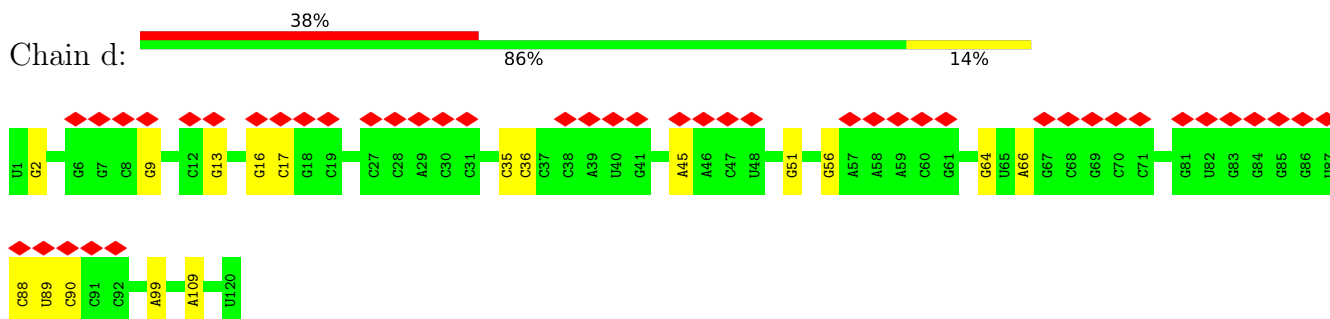
• Molecule 39: 50S ribosomal protein L27



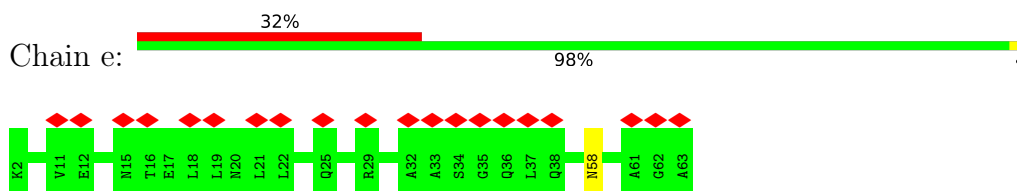
• Molecule 40: 50S ribosomal protein L28



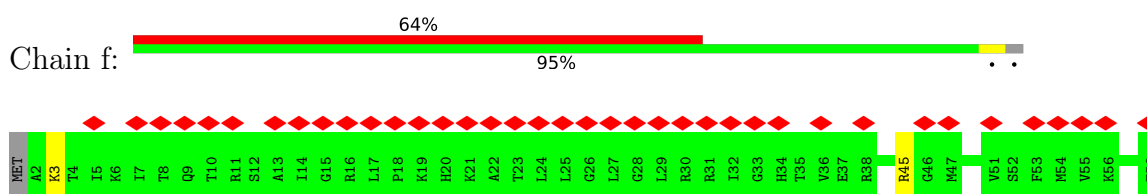
- Molecule 41: 5S rRNA



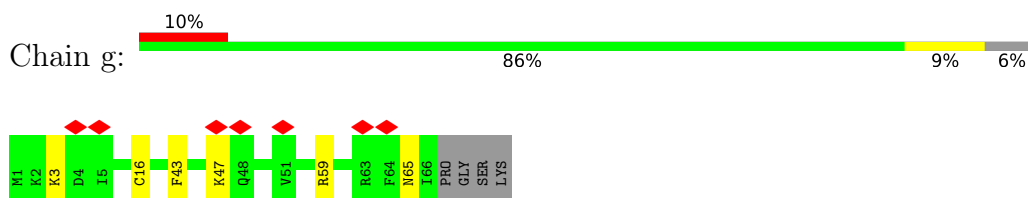
- Molecule 42: 50S ribosomal protein L29



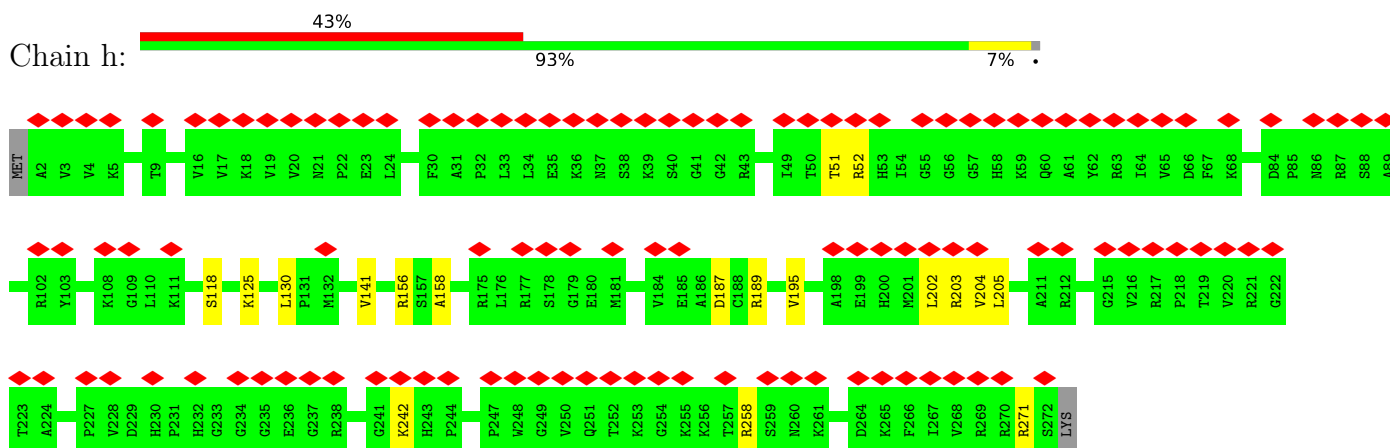
- Molecule 43: 50S ribosomal protein L30



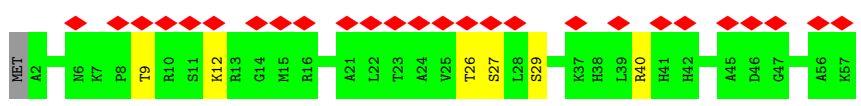
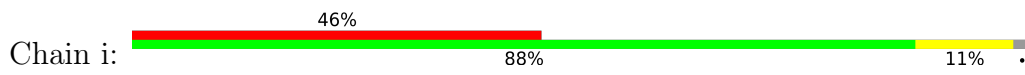
- Molecule 44: 50S ribosomal protein L31



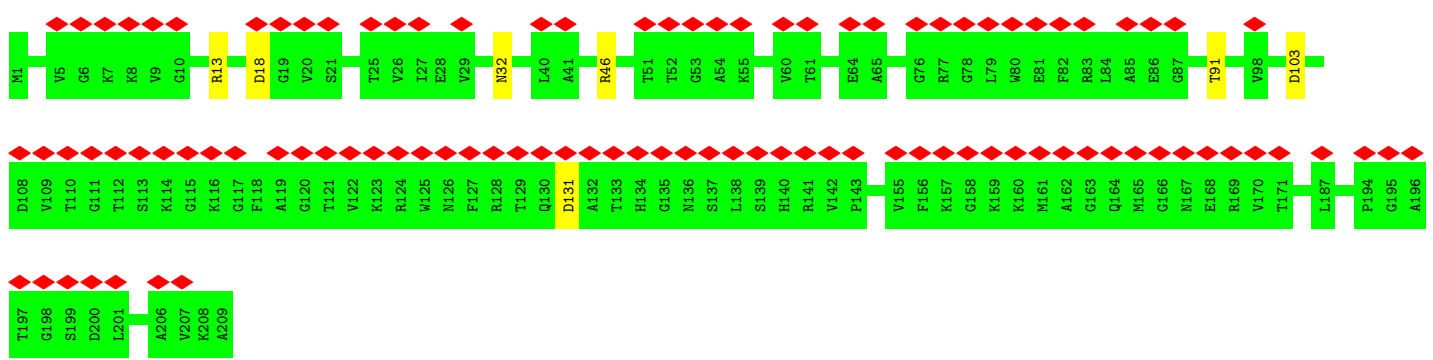
- Molecule 45: 50S ribosomal protein L2



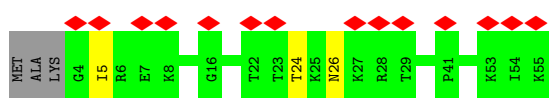
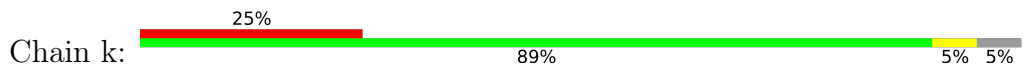
- Molecule 46: 50S ribosomal protein L32



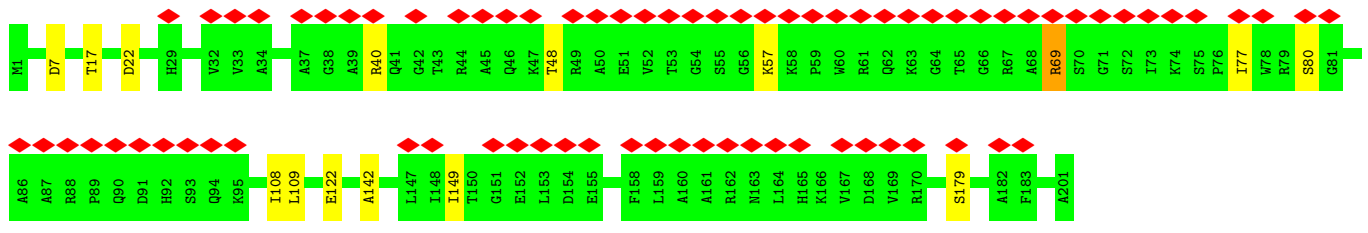
• Molecule 47: 50S ribosomal protein L3



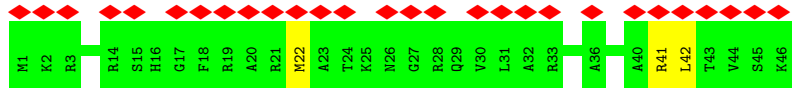
• Molecule 48: 50S ribosomal protein L33



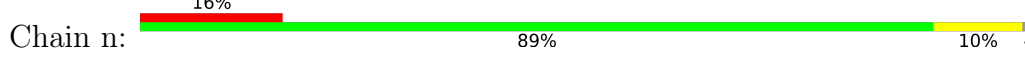
• Molecule 49: 50S ribosomal protein L4

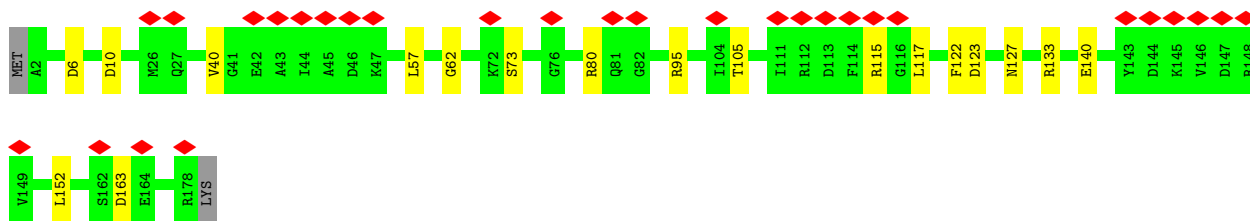


• Molecule 50: 50S ribosomal protein L34

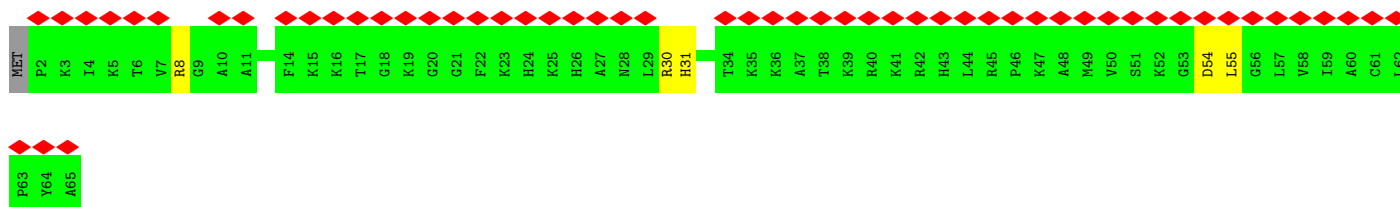
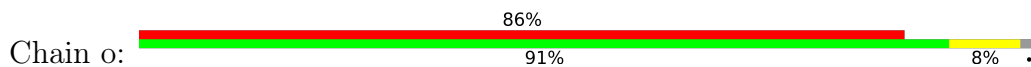


• Molecule 51: 50S ribosomal protein L5

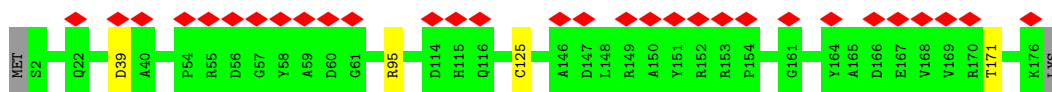




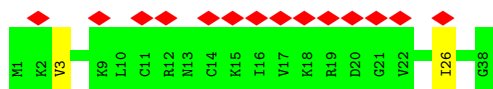
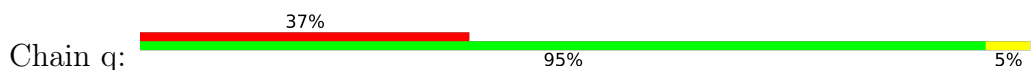
- Molecule 52: 50S ribosomal protein L35



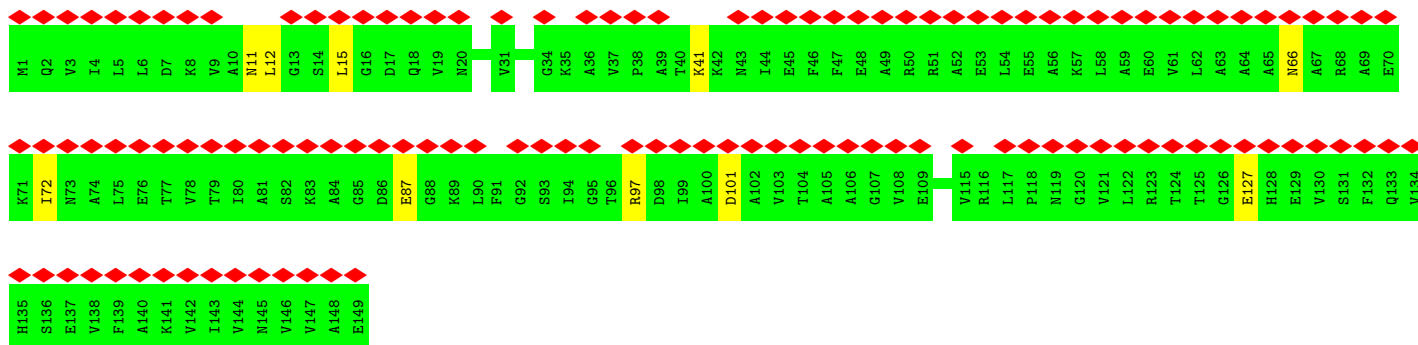
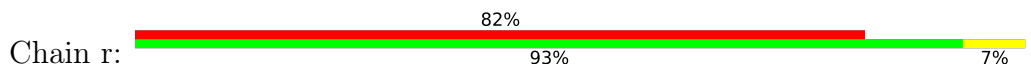
- Molecule 53: 50S ribosomal protein L6



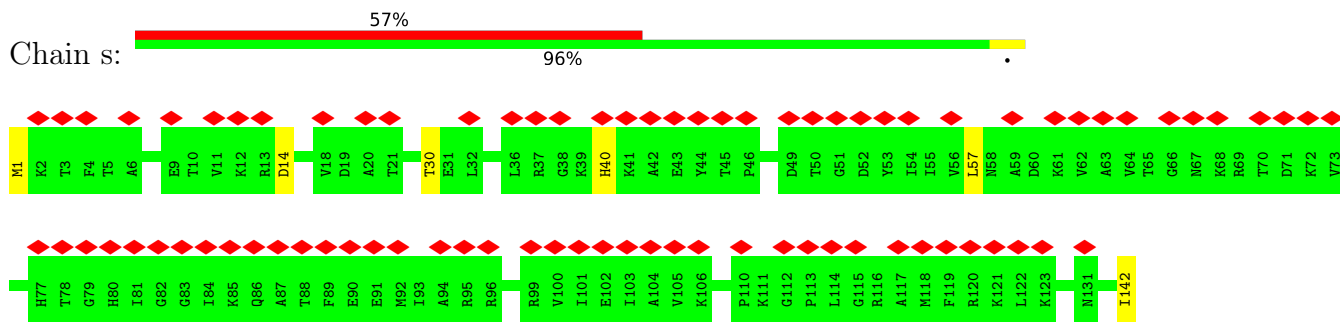
- Molecule 54: 50S ribosomal protein L36



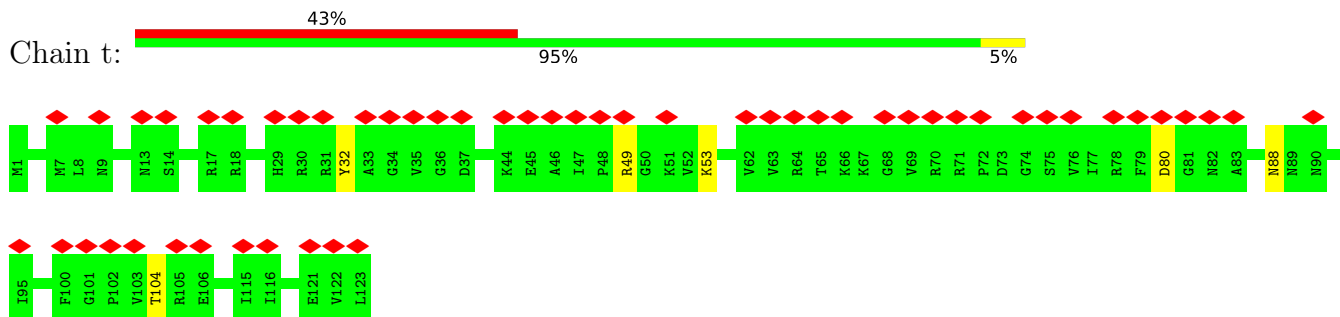
- Molecule 55: 50S ribosomal protein L9



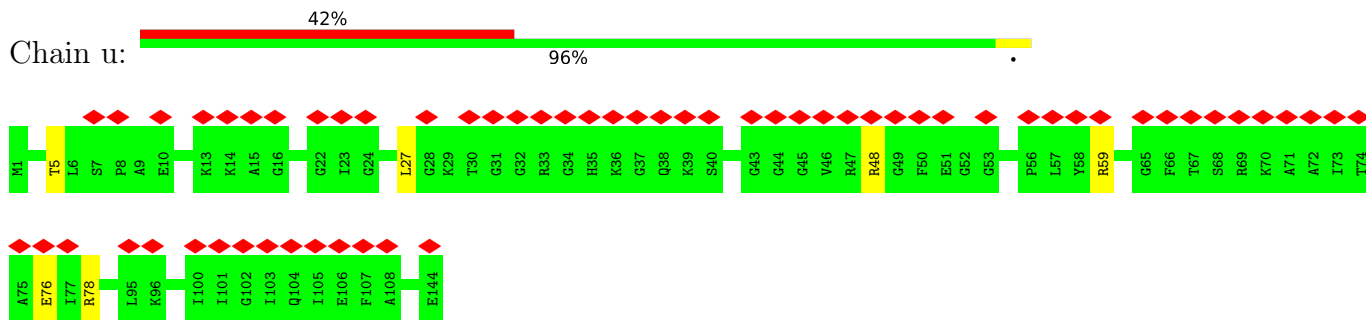
- Molecule 56: 50S ribosomal protein L13



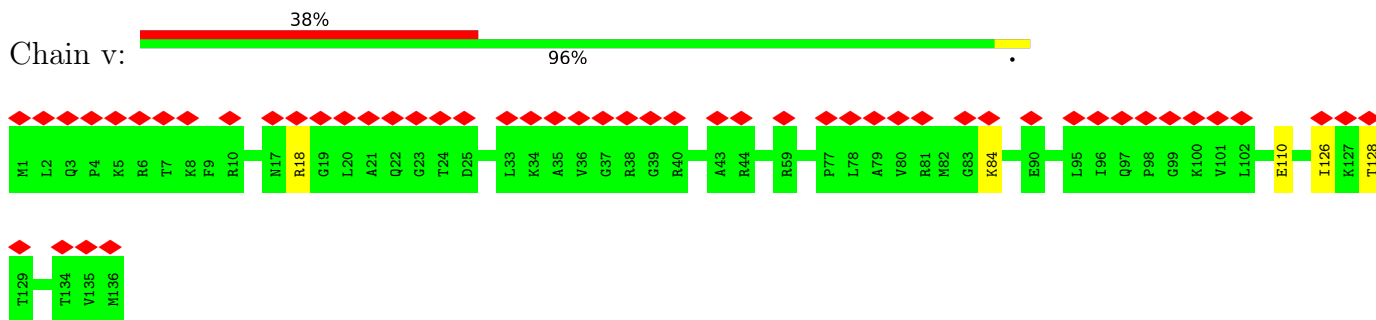
• Molecule 57: 50S ribosomal protein L14



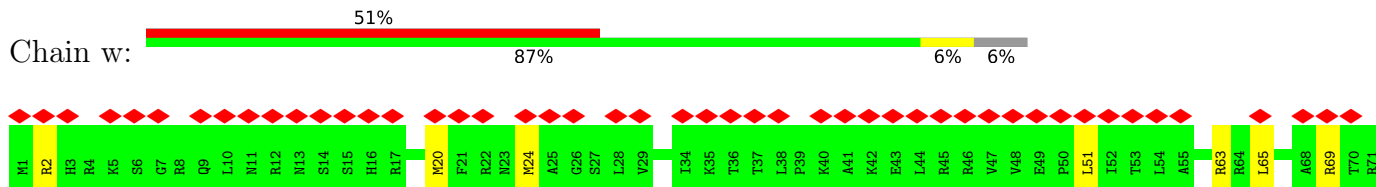
• Molecule 58: 50S ribosomal protein L15



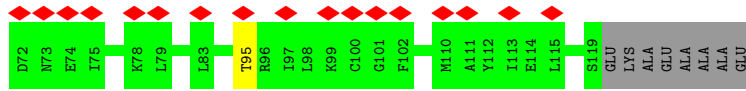
• Molecule 59: 50S ribosomal protein L16



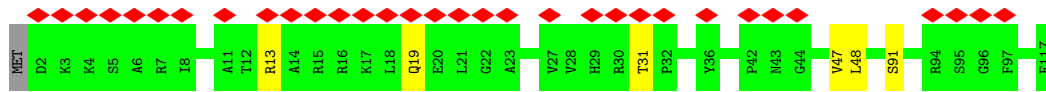
• Molecule 60: 50S ribosomal protein L17



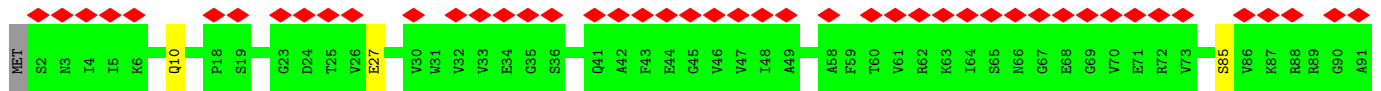
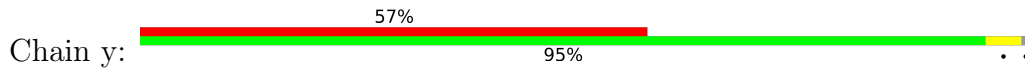




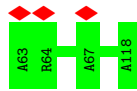
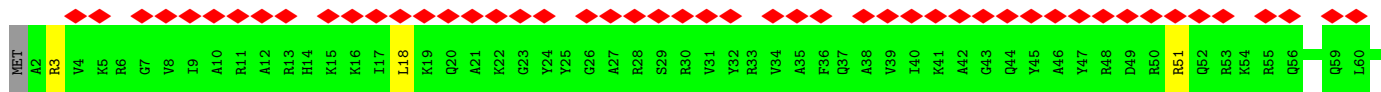
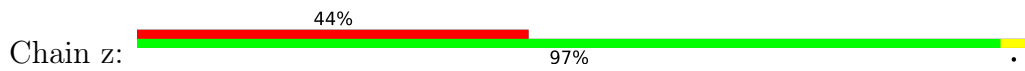
• Molecule 61: 50S ribosomal protein L18



• Molecule 62: 50S ribosomal protein L19



• Molecule 63: 50S ribosomal protein L20



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	5913	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 ( $e^-/\text{\AA}^2$ )	45	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.204	Depositor
Minimum map value	-0.113	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.017	Depositor
Map size (Å)	564.4608, 564.4608, 564.4608	wwPDB
Map dimensions	192, 192, 192	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	2.9399002, 2.9399002, 2.9399002	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	0	0.37	0/829	0.67	0/1107
2	1	0.48	0/864	0.82	0/1156
3	2	0.41	0/752	0.71	0/1005
4	3	0.35	0/796	0.67	2/1062 (0.2%)
5	4	0.40	0/766	0.68	0/1025
6	5	1.13	6/528 (1.1%)	0.97	1/810 (0.1%)
7	6	1.12	4/603 (0.7%)	0.97	0/926
8	7	0.96	4/388 (1.0%)	1.04	0/604
9	A	0.38	0/1810	0.75	1/2821 (0.0%)
9	B	0.43	0/1810	0.85	7/2821 (0.2%)
10	AA	0.58	2/9768 (0.0%)	0.74	11/13191 (0.1%)
11	AB	0.43	0/808	0.60	0/1088
12	AC	0.47	0/1808	0.61	1/2450 (0.0%)
12	AD	0.39	0/1789	0.56	0/2425
13	AE	0.52	4/10545 (0.0%)	0.66	6/14236 (0.0%)
14	AF	0.47	0/657	0.67	0/886
15	C	0.48	0/553	0.82	0/743
16	D	0.31	4/36557 (0.0%)	0.72	20/57006 (0.0%)
17	E	0.57	0/675	0.85	0/895
18	F	0.56	0/597	0.87	0/792
19	G	0.48	0/1791	0.71	0/2413
20	H	0.54	1/1746 (0.1%)	1.03	12/2382 (0.5%)
21	I	0.43	0/1663	0.71	0/2241
22	J	0.46	0/1665	0.74	0/2227
23	K	0.45	0/1165	0.75	0/1568
24	L	0.43	0/867	0.75	1/1171 (0.1%)
25	M	0.50	0/1195	0.81	0/1602
26	N	0.41	0/989	0.69	0/1326
27	O	0.43	0/1034	0.75	0/1375
28	P	0.39	0/800	0.75	0/1082
29	Q	0.40	0/893	0.69	0/1205
30	R	0.35	0/952	0.74	0/1274

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
31	S	0.49	0/817	0.78	0/1088
32	T	0.53	0/722	0.86	0/964
33	U	0.44	0/659	0.78	0/884
34	V	0.34	0/657	0.62	0/881
35	W	0.38	0/680	0.62	0/915
36	X	0.48	0/909	0.87	0/1215
37	Y	0.26	0/65	0.74	0/98
38	a	0.37	3/69247 (0.0%)	0.72	16/107985 (0.0%)
39	b	0.39	0/589	0.71	0/779
40	c	0.48	0/635	0.81	1/848 (0.1%)
41	d	0.28	0/2872	0.69	0/4478
42	e	0.53	0/502	0.82	0/667
43	f	0.44	0/452	0.78	0/605
44	g	0.43	0/531	0.69	0/709
45	h	0.39	0/2121	0.78	0/2852
46	i	0.40	0/450	0.79	0/599
47	j	0.43	0/1586	0.70	0/2134
48	k	0.34	0/433	0.65	0/576
49	l	0.46	0/1571	0.77	2/2113 (0.1%)
50	m	0.53	0/380	0.99	0/498
51	n	0.49	0/1434	0.88	3/1926 (0.2%)
52	o	0.45	0/513	0.83	0/676
53	p	0.39	0/1333	0.67	0/1805
54	q	0.37	0/303	0.77	0/397
55	r	0.43	0/1122	0.69	0/1515
56	s	0.49	0/1152	0.75	0/1551
57	t	0.41	0/955	0.77	0/1279
58	u	0.40	0/1062	0.76	0/1413
59	v	0.46	0/1093	0.81	0/1460
60	w	0.52	0/964	0.87	0/1289
61	x	0.46	0/902	0.81	0/1209
62	y	0.41	0/929	0.72	1/1242 (0.1%)
63	z	0.60	0/960	0.91	0/1278
All	All	0.42	28/186263 (0.0%)	0.73	85/274843 (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
9	A	0	2

*Continued on next page...*

Continued from previous page...

Mol	Chain	#Chirality outliers	#Planarity outliers
9	B	0	2
10	AA	0	11
13	AE	0	5
14	AF	0	1
20	H	0	3
36	X	0	1
All	All	0	25

The worst 5 of 28 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	D	1339	A	O3'-P	10.49	1.73	1.61
10	AA	374	GLU	C-N	10.35	1.53	1.34
13	AE	88	CYS	CB-SG	-10.16	1.65	1.82
6	5	109	DT	O3'-P	8.55	1.71	1.61
7	6	10	DG	C1'-N9	-8.31	1.35	1.47

The worst 5 of 85 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	AA	1250	SER	C-N-CA	11.19	149.67	121.70
38	a	2252	G	N9-C1'-C2'	-10.99	99.72	114.00
16	D	1401	G	N9-C1'-C2'	-10.72	100.07	114.00
51	n	73	SER	N-CA-CB	-10.61	94.59	110.50
16	D	1499	A	N9-C1'-C2'	-10.33	100.57	114.00

There are no chirality outliers.

5 of 25 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
9	A	19	G	Sidechain
9	A	7	G	Sidechain
10	AA	205	PRO	Peptide
10	AA	594	VAL	Peptide
10	AA	595	THR	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

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
1	0	101/103 (98%)	97 (96%)	3 (3%)	1 (1%)	15	55
2	1	108/110 (98%)	104 (96%)	4 (4%)	0	100	100
3	2	92/100 (92%)	90 (98%)	2 (2%)	0	100	100
4	3	101/104 (97%)	96 (95%)	4 (4%)	1 (1%)	15	55
5	4	92/94 (98%)	90 (98%)	2 (2%)	0	100	100
10	AA	1217/1342 (91%)	1071 (88%)	124 (10%)	22 (2%)	8	40
11	AB	94/181 (52%)	88 (94%)	6 (6%)	0	100	100
12	AC	228/329 (69%)	214 (94%)	12 (5%)	2 (1%)	17	57
12	AD	226/329 (69%)	211 (93%)	15 (7%)	0	100	100
13	AE	1329/1407 (94%)	1199 (90%)	121 (9%)	9 (1%)	22	63
14	AF	81/91 (89%)	74 (91%)	7 (9%)	0	100	100
15	C	64/75 (85%)	63 (98%)	1 (2%)	0	100	100
17	E	84/87 (97%)	83 (99%)	1 (1%)	0	100	100
18	F	68/71 (96%)	68 (100%)	0	0	100	100
19	G	223/241 (92%)	210 (94%)	13 (6%)	0	100	100
20	H	255/557 (46%)	189 (74%)	54 (21%)	12 (5%)	2	21
21	I	206/208 (99%)	196 (95%)	9 (4%)	1 (0%)	29	69
22	J	203/206 (98%)	198 (98%)	5 (2%)	0	100	100
23	K	154/167 (92%)	146 (95%)	7 (4%)	1 (1%)	25	66
24	L	102/135 (76%)	97 (95%)	4 (4%)	1 (1%)	15	55
25	M	149/179 (83%)	144 (97%)	4 (3%)	1 (1%)	22	63
26	N	127/130 (98%)	121 (95%)	5 (4%)	1 (1%)	19	60
27	O	125/130 (96%)	115 (92%)	9 (7%)	1 (1%)	19	60
28	P	97/103 (94%)	89 (92%)	7 (7%)	1 (1%)	15	55
29	Q	115/129 (89%)	104 (90%)	9 (8%)	2 (2%)	9	42

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
30	R	117/124 (94%)	116 (99%)	1 (1%)	0	100	100
31	S	98/101 (97%)	96 (98%)	2 (2%)	0	100	100
32	T	86/89 (97%)	82 (95%)	4 (5%)	0	100	100
33	U	80/82 (98%)	75 (94%)	4 (5%)	1 (1%)	12	48
34	V	78/84 (93%)	74 (95%)	4 (5%)	0	100	100
35	W	81/92 (88%)	78 (96%)	3 (4%)	0	100	100
36	X	114/118 (97%)	107 (94%)	5 (4%)	2 (2%)	8	40
39	b	74/85 (87%)	69 (93%)	5 (7%)	0	100	100
40	c	75/78 (96%)	72 (96%)	3 (4%)	0	100	100
42	e	60/62 (97%)	57 (95%)	3 (5%)	0	100	100
43	f	56/59 (95%)	53 (95%)	3 (5%)	0	100	100
44	g	64/70 (91%)	63 (98%)	1 (2%)	0	100	100
45	h	269/273 (98%)	259 (96%)	9 (3%)	1 (0%)	34	72
46	i	54/57 (95%)	51 (94%)	3 (6%)	0	100	100
47	j	207/209 (99%)	198 (96%)	9 (4%)	0	100	100
48	k	50/55 (91%)	50 (100%)	0	0	100	100
49	l	199/201 (99%)	190 (96%)	8 (4%)	1 (0%)	29	69
50	m	44/46 (96%)	43 (98%)	1 (2%)	0	100	100
51	n	175/179 (98%)	162 (93%)	11 (6%)	2 (1%)	14	52
52	o	62/65 (95%)	59 (95%)	3 (5%)	0	100	100
53	p	173/177 (98%)	161 (93%)	12 (7%)	0	100	100
54	q	36/38 (95%)	35 (97%)	1 (3%)	0	100	100
55	r	147/149 (99%)	136 (92%)	11 (8%)	0	100	100
56	s	140/142 (99%)	135 (96%)	5 (4%)	0	100	100
57	t	121/123 (98%)	111 (92%)	10 (8%)	0	100	100
58	u	142/144 (99%)	135 (95%)	7 (5%)	0	100	100
59	v	134/136 (98%)	129 (96%)	5 (4%)	0	100	100
60	w	117/127 (92%)	107 (92%)	10 (8%)	0	100	100
61	x	114/117 (97%)	108 (95%)	6 (5%)	0	100	100
62	y	112/115 (97%)	105 (94%)	7 (6%)	0	100	100
63	z	115/118 (98%)	110 (96%)	4 (4%)	1 (1%)	17	57

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	9035/10123 (89%)	8383 (93%)	588 (6%)	64 (1%)	26	63

5 of 64 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	AA	596	ASP
10	AA	853	ASP
10	AA	859	GLU
10	AA	862	LEU
10	AA	937	ASP

### 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	
1	0	84/84 (100%)	78 (93%)	6 (7%)	14	39
2	1	93/93 (100%)	84 (90%)	9 (10%)	8	27
3	2	81/84 (96%)	76 (94%)	5 (6%)	18	43
4	3	84/85 (99%)	78 (93%)	6 (7%)	14	39
5	4	78/78 (100%)	74 (95%)	4 (5%)	24	48
10	AA	1054/1157 (91%)	1011 (96%)	43 (4%)	30	55
11	AB	86/158 (54%)	84 (98%)	2 (2%)	50	70
12	AC	198/286 (69%)	184 (93%)	14 (7%)	14	39
12	AD	196/286 (68%)	194 (99%)	2 (1%)	76	86
13	AE	1120/1168 (96%)	1051 (94%)	69 (6%)	18	43
14	AF	70/75 (93%)	70 (100%)	0	100	100
15	C	57/65 (88%)	55 (96%)	2 (4%)	36	59
17	E	65/66 (98%)	60 (92%)	5 (8%)	13	37
18	F	60/61 (98%)	57 (95%)	3 (5%)	24	49
19	G	187/199 (94%)	179 (96%)	8 (4%)	29	53
20	H	137/461 (30%)	128 (93%)	9 (7%)	16	41

Continued on next page...



*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
21	I	171/171 (100%)	166 (97%)	5 (3%)	42	64
22	J	172/173 (99%)	166 (96%)	6 (4%)	36	59
23	K	119/126 (94%)	112 (94%)	7 (6%)	19	45
24	L	91/116 (78%)	85 (93%)	6 (7%)	16	41
25	M	124/147 (84%)	116 (94%)	8 (6%)	17	42
26	N	104/105 (99%)	102 (98%)	2 (2%)	57	75
27	O	105/107 (98%)	100 (95%)	5 (5%)	25	51
28	P	86/90 (96%)	80 (93%)	6 (7%)	15	40
29	Q	90/99 (91%)	87 (97%)	3 (3%)	38	61
30	R	101/104 (97%)	94 (93%)	7 (7%)	15	40
31	S	83/84 (99%)	79 (95%)	4 (5%)	25	51
32	T	76/77 (99%)	64 (84%)	12 (16%)	2	13
33	U	65/65 (100%)	60 (92%)	5 (8%)	13	37
34	V	74/78 (95%)	72 (97%)	2 (3%)	44	65
35	W	72/79 (91%)	68 (94%)	4 (6%)	21	46
36	X	94/96 (98%)	85 (90%)	9 (10%)	8	27
39	b	58/63 (92%)	57 (98%)	1 (2%)	60	78
40	c	67/68 (98%)	64 (96%)	3 (4%)	27	52
42	e	54/54 (100%)	53 (98%)	1 (2%)	57	75
43	f	48/49 (98%)	46 (96%)	2 (4%)	30	54
44	g	59/62 (95%)	53 (90%)	6 (10%)	7	25
45	h	216/218 (99%)	199 (92%)	17 (8%)	12	35
46	i	47/48 (98%)	41 (87%)	6 (13%)	4	18
47	j	164/164 (100%)	157 (96%)	7 (4%)	29	53
48	k	47/49 (96%)	44 (94%)	3 (6%)	17	42
49	l	165/165 (100%)	151 (92%)	14 (8%)	10	33
50	m	38/38 (100%)	35 (92%)	3 (8%)	12	35
51	n	148/150 (99%)	134 (90%)	14 (10%)	8	27
52	o	51/52 (98%)	46 (90%)	5 (10%)	8	26
53	p	136/138 (99%)	132 (97%)	4 (3%)	42	64
54	q	34/34 (100%)	32 (94%)	2 (6%)	19	45

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
55	r	114/114 (100%)	104 (91%)	10 (9%)	10	31
56	s	116/116 (100%)	110 (95%)	6 (5%)	23	48
57	t	104/104 (100%)	98 (94%)	6 (6%)	20	45
58	u	103/103 (100%)	97 (94%)	6 (6%)	20	45
59	v	109/109 (100%)	104 (95%)	5 (5%)	27	52
60	w	99/103 (96%)	91 (92%)	8 (8%)	11	35
61	x	86/87 (99%)	80 (93%)	6 (7%)	15	40
62	y	99/100 (99%)	95 (96%)	4 (4%)	31	55
63	z	89/90 (99%)	87 (98%)	2 (2%)	52	71
All	All	7528/8401 (90%)	7109 (94%)	419 (6%)	25	46

5 of 419 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
28	P	37	ARG
44	g	16	CYS
60	w	2	ARG
30	R	56	ARG
33	U	6	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
19	G	18	HIS
23	K	70	ASN
36	X	105	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
16	D	1513/1542 (98%)	284 (18%)	33 (2%)
37	Y	2/3 (66%)	2 (100%)	0
38	a	2859/2904 (98%)	533 (18%)	0
41	d	119/120 (99%)	17 (14%)	0
8	7	15/38 (39%)	7 (46%)	0
9	A	75/76 (98%)	29 (38%)	6 (8%)
9	B	75/76 (98%)	34 (45%)	5 (6%)

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	4658/4759 (97%)	906 (19%)	44 (0%)

5 of 906 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
8	7	56	U
8	7	57	G
8	7	58	A
8	7	59	U
8	7	60	U

5 of 44 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
16	D	793	U
16	D	1212	U
16	D	991	U
16	D	1145	A
16	D	1214	C

## 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 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

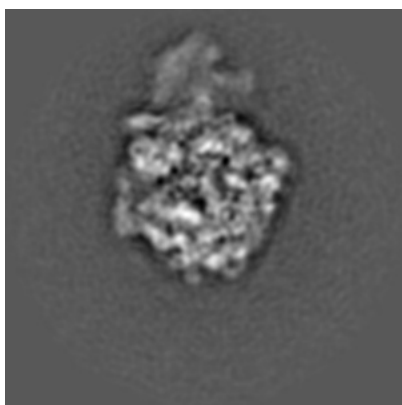
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21477. These allow visual inspection of the internal detail of the map and identification of artifacts.

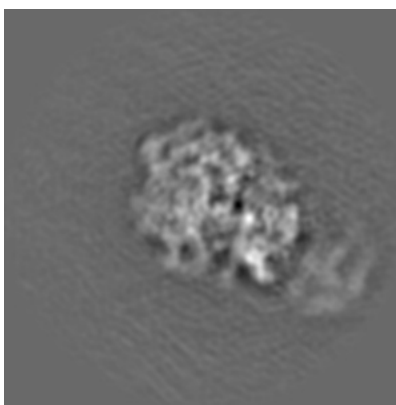
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

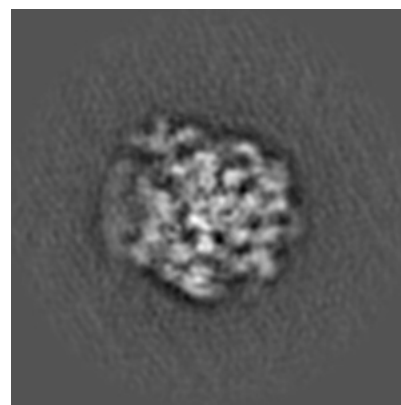
#### 6.1.1 Primary map



X



Y

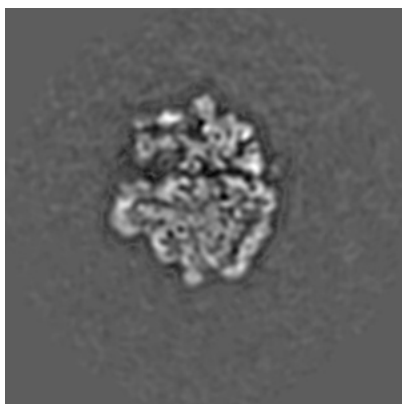


Z

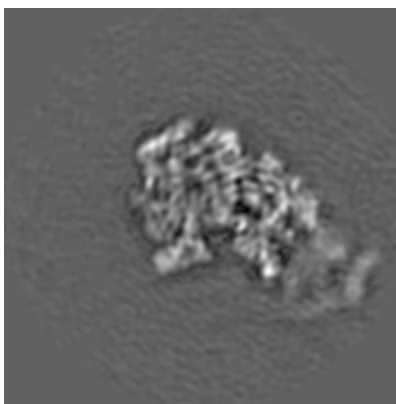
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

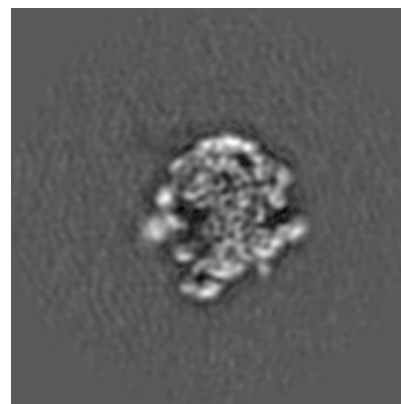
#### 6.2.1 Primary map



X Index: 96



Y Index: 96

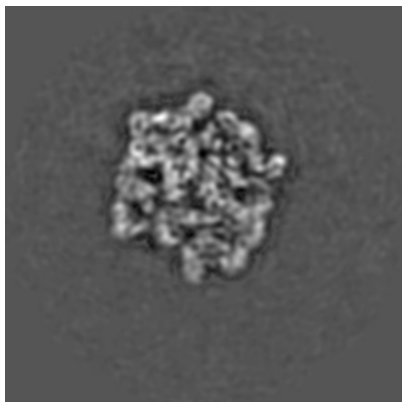


Z Index: 96

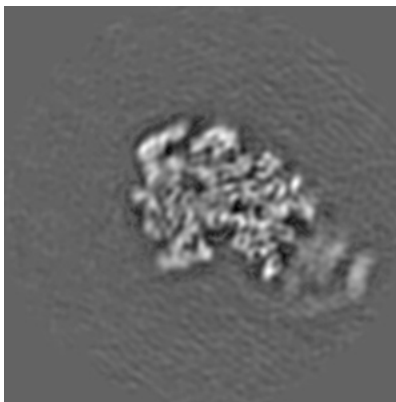
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

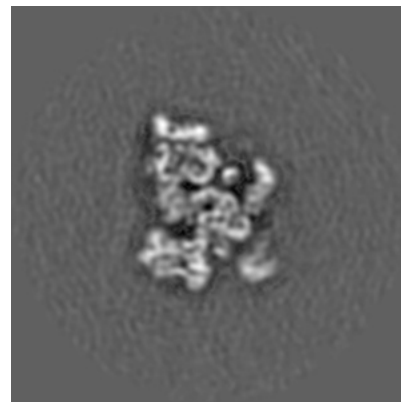
### 6.3.1 Primary map



X Index: 91



Y Index: 98

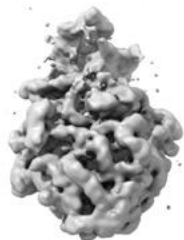


Z Index: 117

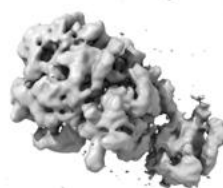
The images above show the largest variance slices of the map in three orthogonal directions.

## 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 0.017. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

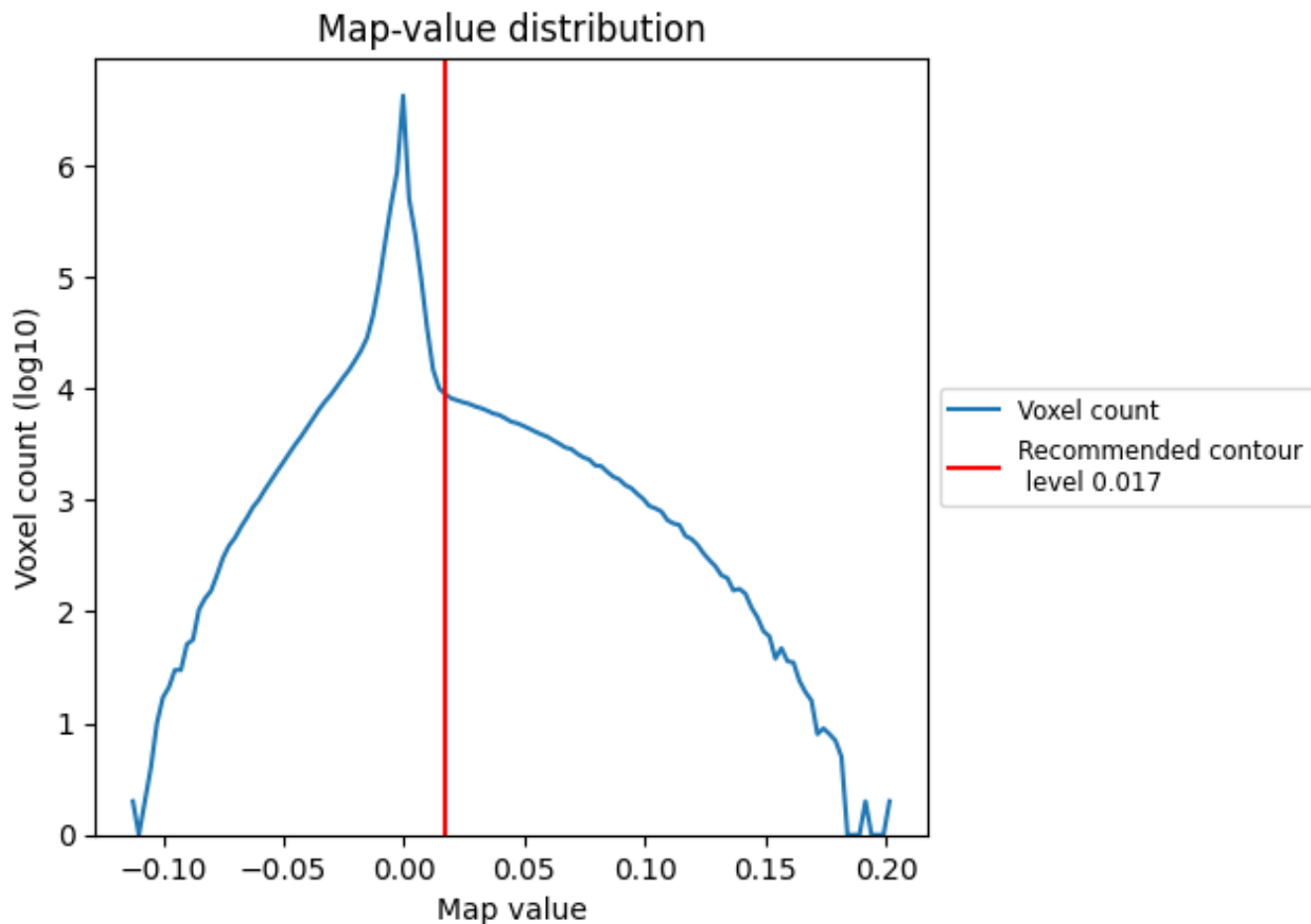
## 6.5 Mask visualisation

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.

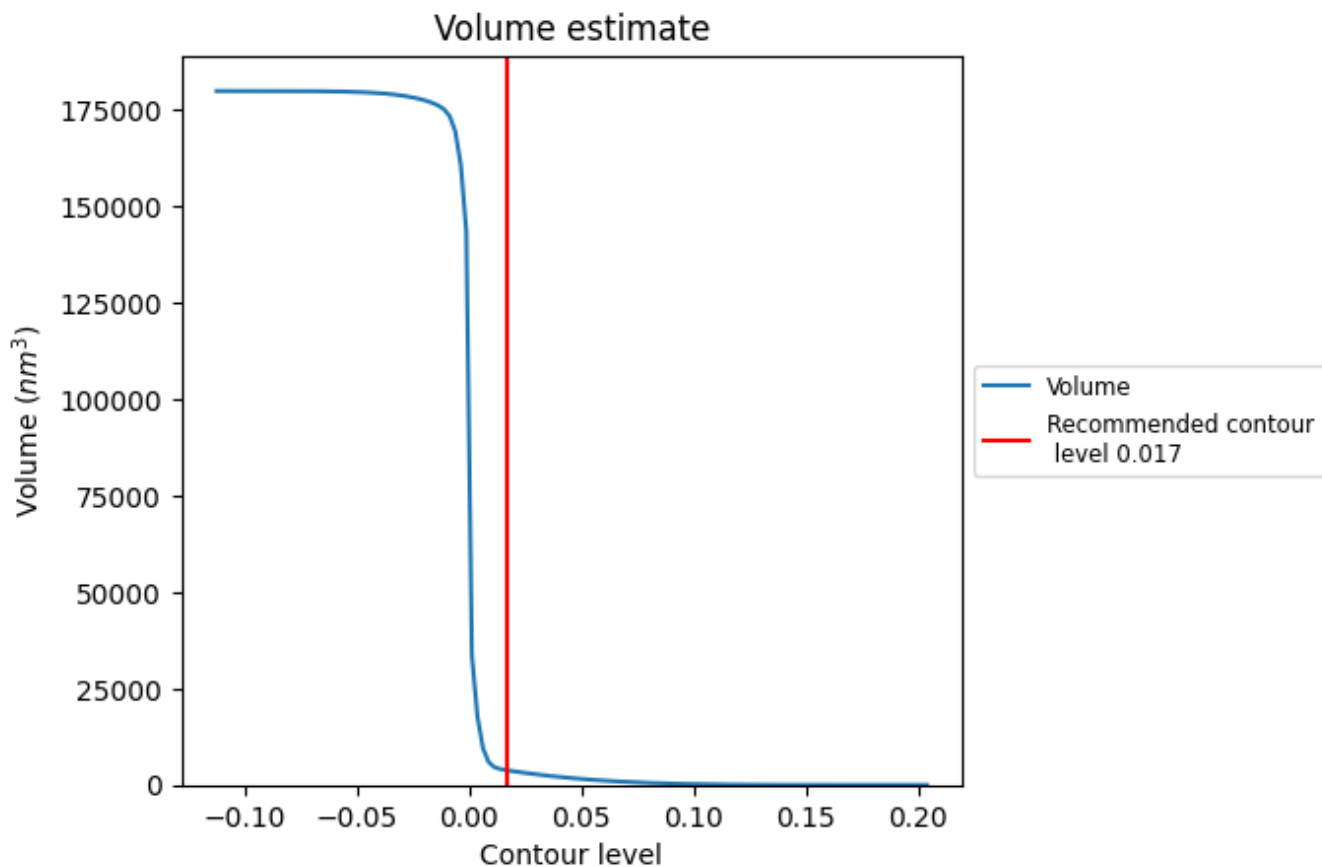
### 7.1 Map-value distribution [i](#)



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.



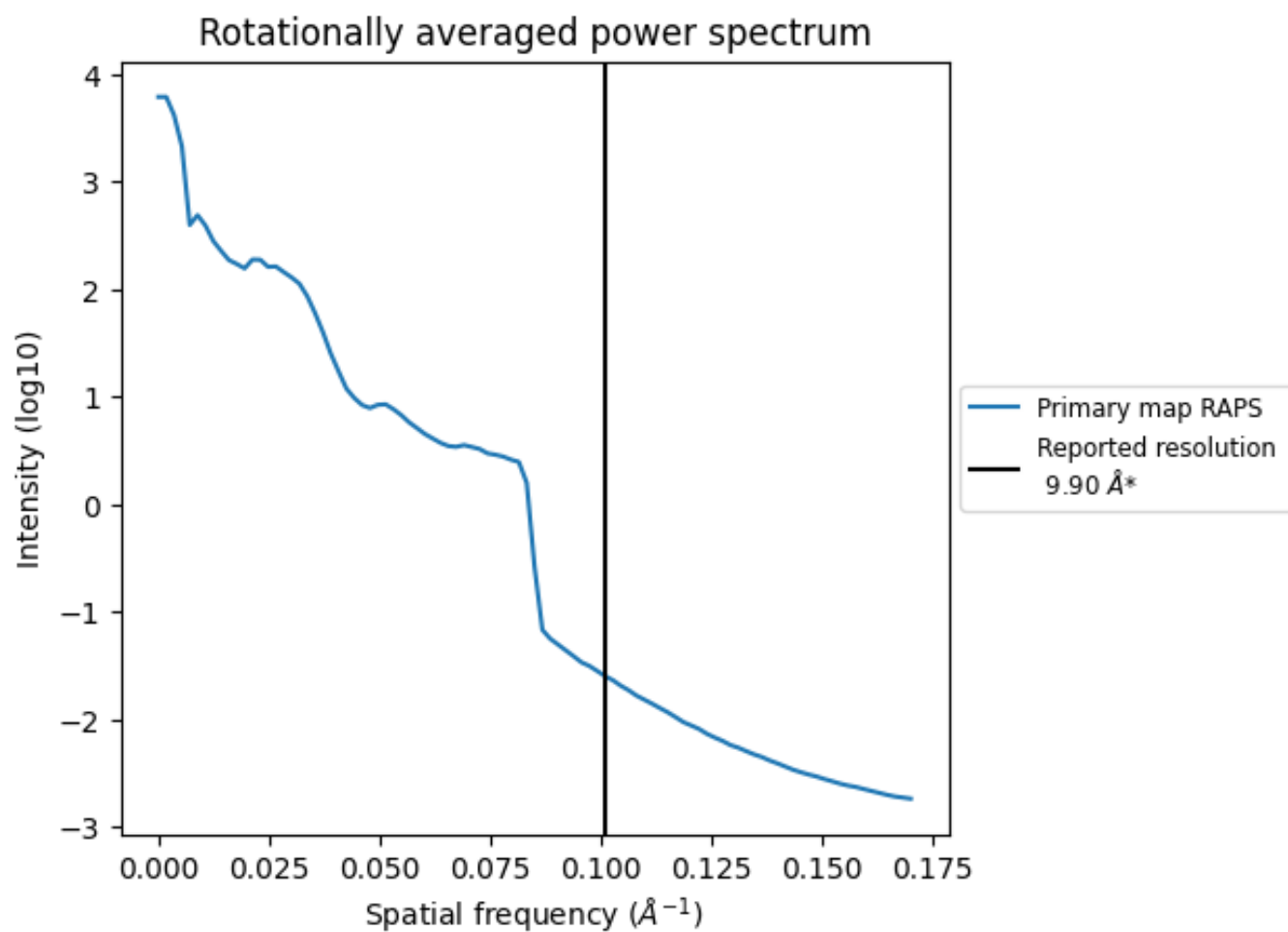
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 3791  $\text{nm}^3$ ; this corresponds to an approximate mass of 3425 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 0.101 Å<sup>-1</sup>

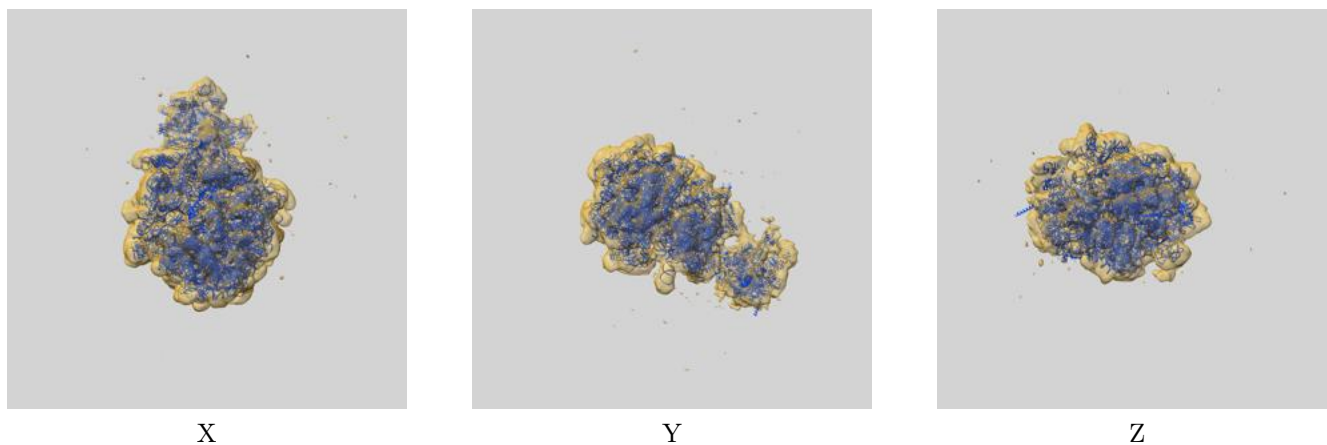
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

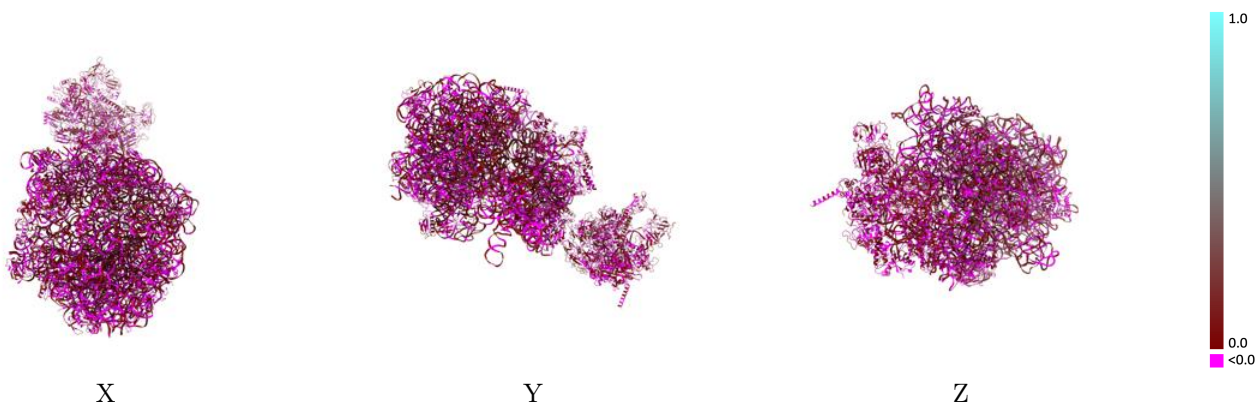
This section contains information regarding the fit between EMDB map EMD-21477 and PDB model 6VYZ. Per-residue inclusion information can be found in section 3 on page 16.

### 9.1 Map-model overlay [i](#)



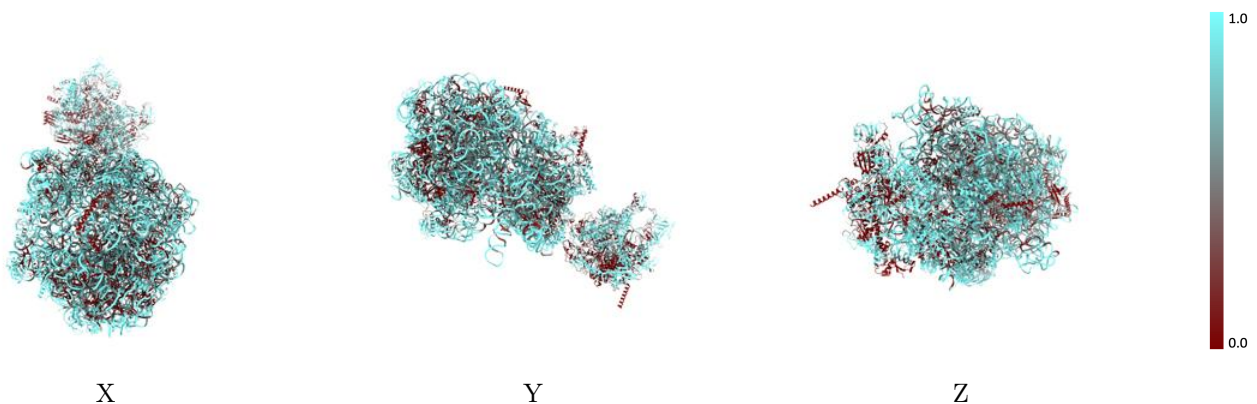
The images above show the 3D surface view of the map at the recommended contour level 0.017 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



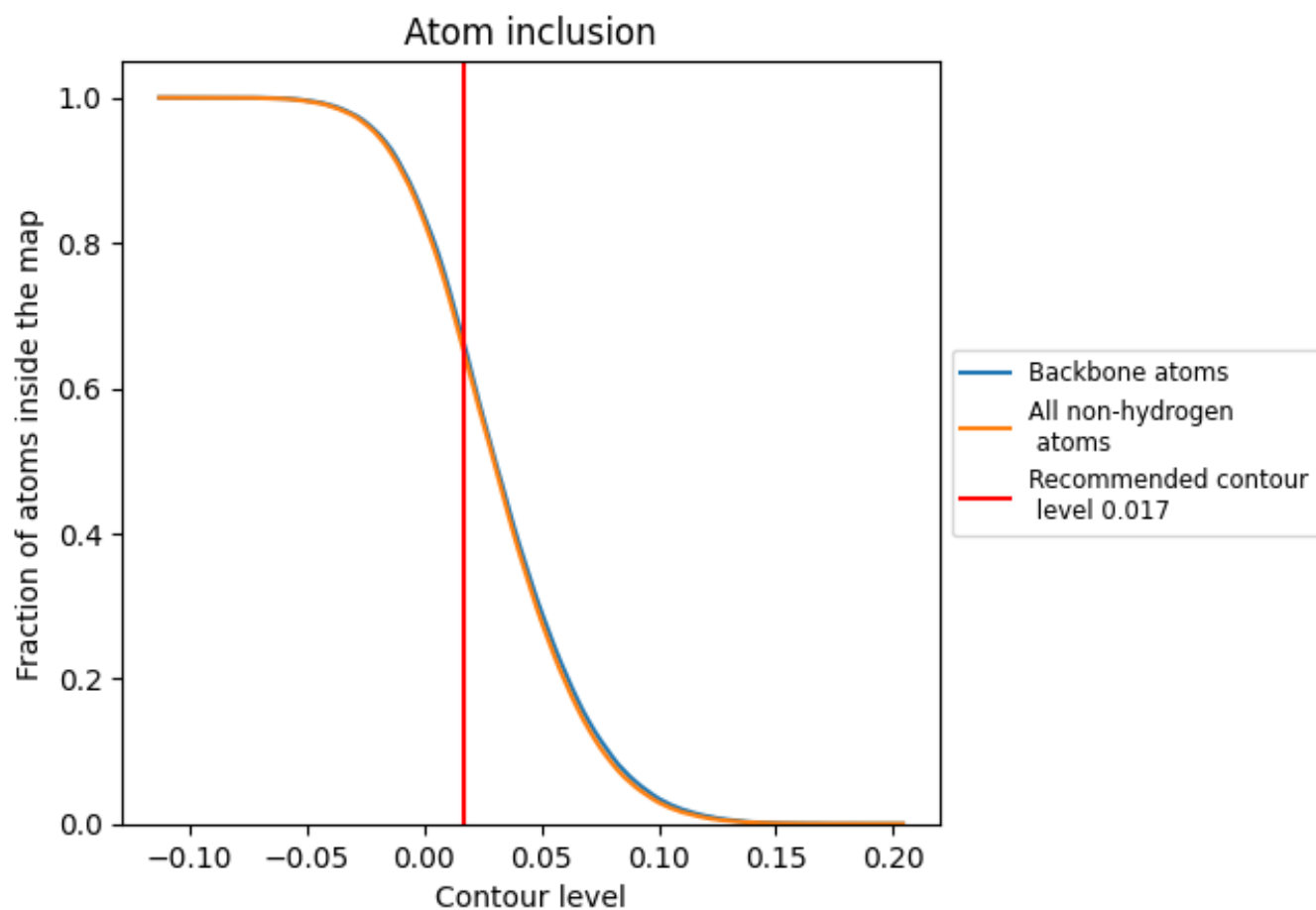
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.017).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 66% of all backbone atoms, 65% 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 (0.017) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6516	 0.0330
0	 0.6926	 0.0040
1	 0.2955	 0.0070
2	 0.5233	 -0.0190
3	 0.7371	 0.0080
4	 0.6978	 0.0180
5	 0.6547	 0.0520
6	 0.6089	 0.0390
7	 0.3718	 0.0540
A	 0.6969	 0.0770
AA	 0.5955	 0.0400
AB	 0.3459	 -0.0030
AC	 0.4794	 0.0480
AD	 0.5020	 0.0210
AE	 0.5282	 0.0260
AF	 0.2692	 -0.0030
B	 0.7475	 0.0610
C	 0.7973	 -0.0040
D	 0.7694	 0.0490
E	 0.4755	 0.0030
F	 0.2361	 0.0230
G	 0.7553	 0.0120
H	 0.3464	 0.0070
I	 0.6000	 0.0040
J	 0.5176	 -0.0110
K	 0.5646	 0.0240
L	 0.8102	 0.0340
M	 0.6460	 0.0070
N	 0.5604	 -0.0260
O	 0.6057	 -0.0190
P	 0.5673	 0.0110
Q	 0.7937	 0.0410
R	 0.6541	 0.0560
S	 0.5142	 -0.0230
T	 0.7899	 0.0220



*Continued on next page...*

Continued from previous page...

Chain	Atom inclusion	Q-score
U	0.4163	-0.0230
V	0.5269	0.0380
W	0.7728	0.0060
X	0.6843	-0.0070
Y	0.0000	-0.0390
a	0.7161	0.0460
b	0.3975	-0.0290
c	0.4476	0.0050
d	0.6216	0.0180
e	0.6810	-0.0050
f	0.3372	-0.0200
g	0.8376	0.0410
h	0.5174	0.0280
i	0.4696	-0.0100
j	0.5124	-0.0170
k	0.6746	0.0190
l	0.5947	0.0100
m	0.3662	0.0140
n	0.7815	0.0190
o	0.1487	-0.0240
p	0.7925	0.0230
q	0.5479	0.0050
r	0.1743	-0.0310
s	0.3827	-0.0210
t	0.5364	0.0210
u	0.5615	0.0090
v	0.5624	0.0340
w	0.3910	-0.0290
x	0.6860	-0.0030
y	0.4313	-0.0270
z	0.5253	-0.0280