



wwPDB EM Validation Summary Report ⓘ

Mar 5, 2026 – 07:24 PM UTC

PDB ID : 8P7Y / pdb_00008p7y
EMDB ID : EMD-17135
Title : Mycoplasma pneumoniae 70S ribosome with second S4 protein on large subunit
Authors : Schacherl, M.; Xue, L.; Spahn, C.M.T.; Mahamid, J.
Deposited on : 2023-05-31
Resolution : 3.70 Å (reported)
Based on initial models : 7OOC, 7OOD

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:

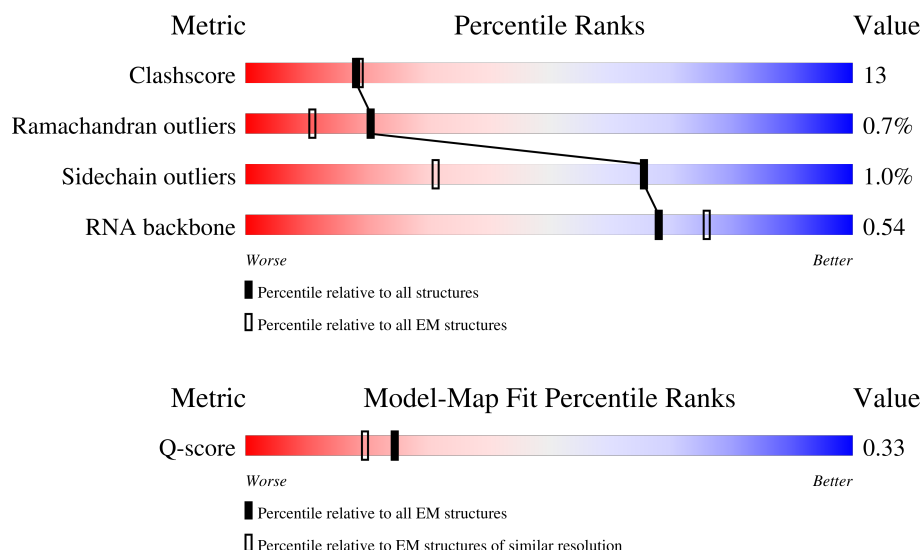
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMD 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

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

The reported resolution of this entry is 3.70 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	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11569 (3.20 - 4.20)

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	0	48	
2	1	59	
3	2	37	

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Mol	Chain	Length	Quality of chain
4	3	2907	
5	4	108	
6	5	1520	
7	6	76	
8	7	75	
9	8	76	
10	A	294	
11	B	273	
12	C	205	
12	U	205	
13	D	219	
14	E	215	
15	F	155	
16	G	142	
17	H	132	
18	I	108	
19	J	121	
20	K	139	
21	L	124	
22	M	61	
23	N	86	
24	O	94	
25	P	85	
26	Q	104	
27	R	87	

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Mol	Chain	Length	Quality of chain
28	S	87	
29	T	60	
30	X	444	
31	Y	21	
32	Z	36	
33	a	287	
34	b	287	
35	c	212	
36	d	180	
37	e	184	
38	f	149	
39	g	161	
40	h	137	
41	i	146	
42	j	122	
43	k	151	
44	l	139	
45	m	124	
46	n	116	
47	o	119	
48	p	127	
49	q	100	
50	r	159	
51	s	237	
52	t	111	

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Mol	Chain	Length	Quality of chain
53	u	104	
54	v	65	
55	w	111	
56	x	97	
57	y	57	
58	z	53	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
65	SPD	5	1601	-	-	X	-

2 Entry composition

There are 67 unique types of molecules in this entry. The entry contains 153255 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 protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	48	Total	C	N	O	S	0	0
			392	242	85	63	2		

- Molecule 2 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	59	Total	C	N	O	S	0	0
			477	300	99	77	1		

- Molecule 3 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	37	Total	C	N	O	S	0	0
			304	189	65	46	4		

- Molecule 4 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	3	2893	Total	C	N	O	P	0	0
			61995	27704	11293	20105	2893		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	4	108	Total	C	N	O	P	0	0
			2305	1030	415	752	108		

- Molecule 6 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	5	1507	Total	C	N	O	P	0	0
			32258	14420	5847	10484	1507		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
5	1003	A	G	conflict	GB 26117688

- Molecule 7 is a RNA chain called tRNA-Ala (E-site).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	6	76	Total	C	N	O	P	0	0
			1620	723	287	534	76		

- Molecule 8 is a RNA chain called tRNA-Asp (P-site).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	7	75	Total	C	N	O	P	0	0
			1599	712	279	533	75		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
7	17	G	-	insertion	GB 26117688
7	55	C	U	conflict	GB 26117688

- Molecule 9 is a RNA chain called tRNA-Lys (A-site).

Mol	Chain	Residues	Atoms					AltConf	Trace
9	8	76	Total	C	N	O	P	0	0
			1615	722	284	533	76		

- Molecule 10 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	A	266	Total	C	N	O	S	0	0
			2138	1359	376	394	9		

- Molecule 11 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	B	232	Total	C	N	O	S	0	0
			1835	1158	343	329	5		

- Molecule 12 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	C	204	Total	C	N	O	S	0	0
			1669	1057	316	292	4		
12	U	204	Total	C	N	O	S	0	0
			1669	1057	316	292	4		

- Molecule 13 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	D	155	Total	C	N	O	S	0	0
			1191	753	228	207	3		

- Molecule 14 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	E	184	Total	C	N	O	S	0	0
			1509	950	270	287	2		

- Molecule 15 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	F	155	Total	C	N	O	S	0	0
			1254	790	240	217	7		

- Molecule 16 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	G	141	Total	C	N	O	S	0	0
			1110	723	193	192	2		

- Molecule 17 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	H	129	Total	C	N	O	S	0	0
			1040	661	195	183	1		

- Molecule 18 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	I	104	Total	C	N	O	S	0	0
			832	536	147	148	1		

- Molecule 19 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	J	114	Total	C	N	O	S	0	0
			829	514	153	156	6		

- Molecule 20 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	K	135	Total	C	N	O	S	0	0
			1071	677	212	180	2		

- Molecule 21 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	L	123	Total	C	N	O		0	0
			991	618	200	173			

- Molecule 22 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	M	60	Total	C	N	O	S	0	0
			474	302	96	72	4		

- Molecule 23 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	N	85	Total	C	N	O		0	0
			689	436	130	123			

- Molecule 24 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	O	87	Total	C	N	O	S	0	0
			705	453	130	118	4		

- Molecule 25 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	P	85	Total	C	N	O	S	0	0
			693	436	138	118	1		

- Molecule 26 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Q	71	Total	C	N	O	S	0	0
			590	378	115	93	4		

- Molecule 27 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	R	86	Total	C	N	O	S	0	0
			700	444	132	122	2		

- Molecule 28 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	S	79	Total	C	N	O		0	0
			643	391	138	114			

- Molecule 29 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	T	59	Total	C	N	O	S	0	0
			519	326	111	80	2		

- Molecule 30 is a protein called Trigger factor.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	X	30	Total	C	N	O	S	0	0
			242	155	43	43	1		

- Molecule 31 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	Y	21	Total	C	N	O	P	0	0
			446	200	80	145	21		

- Molecule 32 is a protein called Nascent chain.

Mol	Chain	Residues	Atoms				AltConf	Trace
32	Z	36	Total	C	N	O	0	0
			187	112	37	38		

- Molecule 33 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	a	285	Total	C	N	O	S	0	0
			2225	1385	437	397	6		

- Molecule 34 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	b	230	Total	C	N	O	S	0	0
			1769	1124	319	319	7		

- Molecule 35 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	c	211	Total	C	N	O	S	0	0
			1654	1053	299	299	3		

- Molecule 36 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	d	179	Total	C	N	O	S	0	0
			1416	910	251	251	4		

- Molecule 37 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms				AltConf	Trace
37	e	176	Total	C	N	O	0	0
			1396	899	247	250		

- Molecule 38 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	f	149	Total	C	N	O	S	0	0
			1210	780	212	215	3		

- Molecule 39 is a protein called 50S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	g	125	Total	C	N	O	S	0	0
			951	606	165	177	3		

- Molecule 40 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	h	128	Total	C	N	O	S	0	0
			959	616	160	177	6		

- Molecule 41 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	i	144	Total	C	N	O	S	0	0
			1164	737	213	209	5		

- Molecule 42 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	j	122	Total	C	N	O	S	0	0
			944	595	178	167	4		

- Molecule 43 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	k	150	Total	C	N	O	S	0	0
			1170	741	228	200	1		

- Molecule 44 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	l	136	Total	C	N	O	S	0	0
			1079	694	196	182	7		

- Molecule 45 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	m	119	Total	C	N	O	S	0	0
			958	609	175	171	3		

- Molecule 46 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	n	116	Total	C	N	O	S	0	0
			918	573	181	162	2		

- Molecule 47 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	o	118	Total	C	N	O	S	0	0
			966	609	186	170	1		

- Molecule 48 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	p	118	Total	C	N	O	S	0	0
			981	624	194	161	2		

- Molecule 49 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	q	99	Total	C	N	O	S	0	0
			811	525	148	134	4		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
50	r	142	Total	C	N	O	S	0	0
			1091	677	212	195	7		

- Molecule 51 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	s	95	Total	C	N	O	S	0	0
			740	486	125	128	1		

- Molecule 52 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	t	111	Total	C	N	O	S	0	0
			871	550	166	152	3		

- Molecule 53 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	u	88	Total	C	N	O	S	0	0
			670	416	132	121	1		

- Molecule 54 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	v	64	Total	C	N	O	S	0	0
			520	320	109	90	1		

- Molecule 55 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	w	110	Total	C	N	O		0	0
			906	576	168	162			

- Molecule 56 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	x	96	Total	C	N	O	S	0	0
			761	481	133	143	4		

- Molecule 57 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	y	56	Total	C	N	O	S	0	0
			452	274	98	75	5		

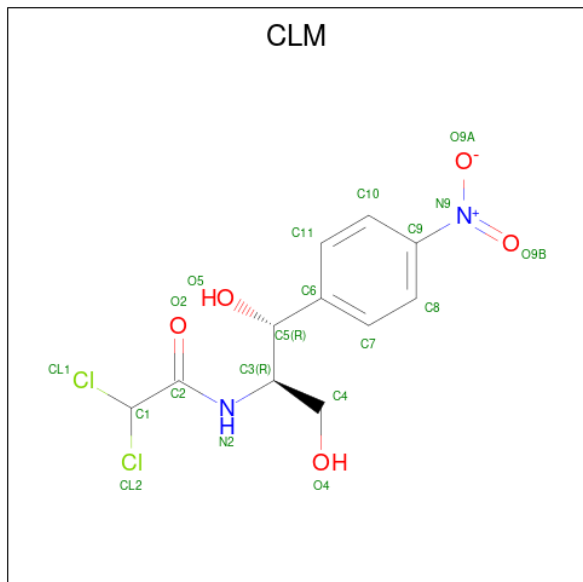
- Molecule 58 is a protein called 50S ribosomal protein L33 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	z	50	Total	C	N	O	S	0	0
			408	255	81	68	4		

- Molecule 59 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
59	2	1	Total	Zn	0
			1	1	
59	M	1	Total	Zn	0
			1	1	
59	Q	1	Total	Zn	0
			1	1	
59	x	1	Total	Zn	0
			1	1	
59	y	1	Total	Zn	0
			1	1	
59	z	1	Total	Zn	0
			1	1	

- Molecule 60 is CHLORAMPHENICOL (CCD ID: CLM) (formula: $C_{11}H_{12}Cl_2N_2O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
60	3	1	Total	C	Cl	N	O	0
			20	11	2	2	5	

- Molecule 61 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
61	3	1	Total	K	0
			1	1	

- Molecule 62 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

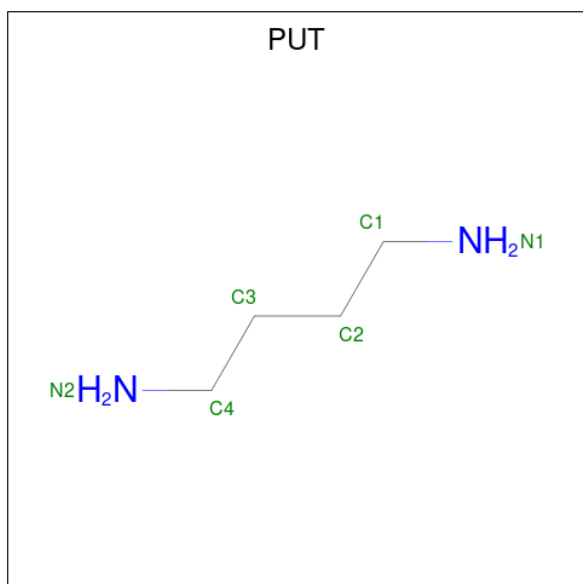
Mol	Chain	Residues	Atoms		AltConf
62	3	221	Total	Mg	0
			221	221	
62	4	1	Total	Mg	0
			1	1	
62	5	90	Total	Mg	0
			90	90	
62	6	1	Total	Mg	0
			1	1	
62	7	2	Total	Mg	0
			2	2	
62	8	2	Total	Mg	0
			2	2	

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Mol	Chain	Residues	Atoms		AltConf
62	K	1	Total 1	Mg 1	0
62	P	1	Total 1	Mg 1	0
62	Y	2	Total 2	Mg 2	0
62	b	2	Total 2	Mg 2	0
62	i	1	Total 1	Mg 1	0
62	y	2	Total 2	Mg 2	0

- Molecule 63 is 1,4-DIAMINOBTANE (CCD ID: PUT) (formula: $C_4H_{12}N_2$).



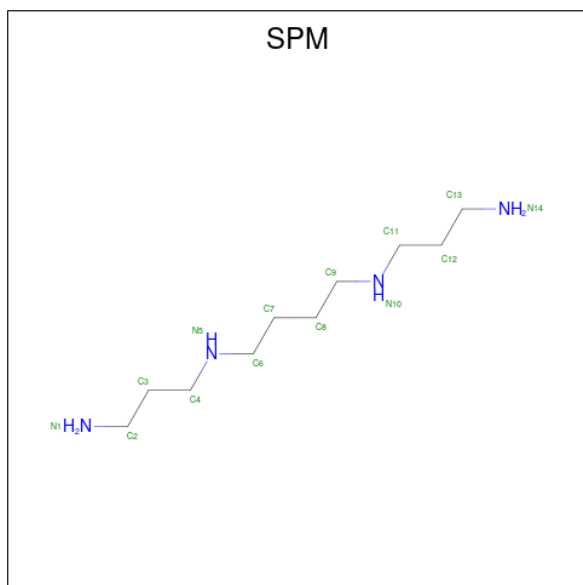
Mol	Chain	Residues	Atoms			AltConf
63	3	1	Total 6	C 4	N 2	0
63	3	1	Total 6	C 4	N 2	0
63	3	1	Total 6	C 4	N 2	0
63	3	1	Total 6	C 4	N 2	0
63	3	1	Total 6	C 4	N 2	0

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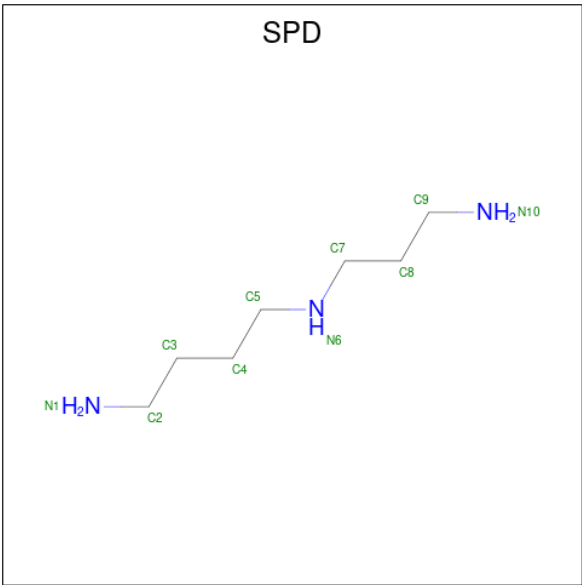
Mol	Chain	Residues	Atoms			AltConf
63	3	1	Total	C	N	0
			6	4	2	
63	3	1	Total	C	N	0
			6	4	2	
63	5	1	Total	C	N	0
			6	4	2	

- Molecule 64 is SPERMINE (CCD ID: SPM) (formula: $C_{10}H_{26}N_4$).



Mol	Chain	Residues	Atoms			AltConf
64	3	1	Total	C	N	0
			14	10	4	
64	3	1	Total	C	N	0
			14	10	4	
64	3	1	Total	C	N	0
			14	10	4	
64	b	1	Total	C	N	0
			14	10	4	

- Molecule 65 is SPERMIDINE (CCD ID: SPD) (formula: $C_7H_{19}N_3$).



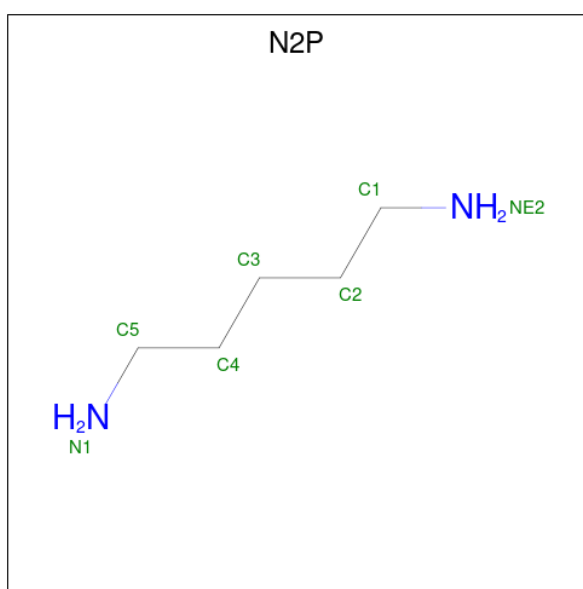
Mol	Chain	Residues	Atoms			AltConf
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	

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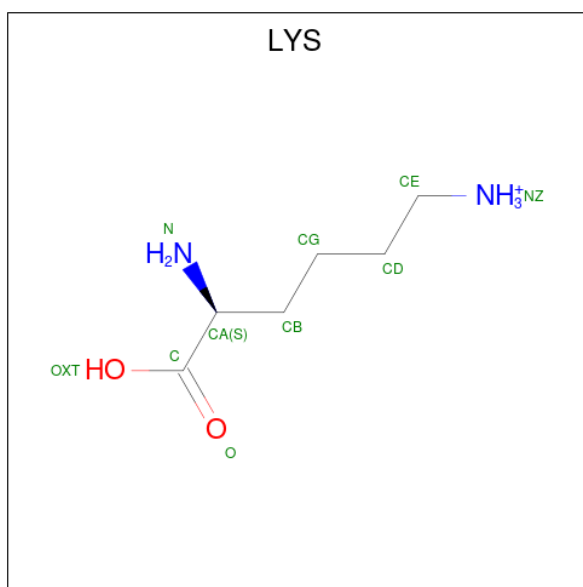
Mol	Chain	Residues	Atoms			AltConf
65	3	1	Total	C	N	0
			10	7	3	
65	3	1	Total	C	N	0
			10	7	3	
65	5	1	Total	C	N	0
			10	7	3	
65	5	1	Total	C	N	0
			10	7	3	

- Molecule 66 is PENTANE-1,5-DIAMINE (CCD ID: N2P) (formula: $C_5H_{14}N_2$).



Mol	Chain	Residues	Atoms			AltConf
66	3	1	Total	C	N	0
			7	5	2	
66	3	1	Total	C	N	0
			7	5	2	
66	3	1	Total	C	N	0
			7	5	2	
66	5	1	Total	C	N	0
			7	5	2	

- Molecule 67 is LYSINE (CCD ID: LYS) (formula: $C_6H_{15}N_2O_2$).

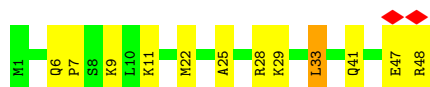
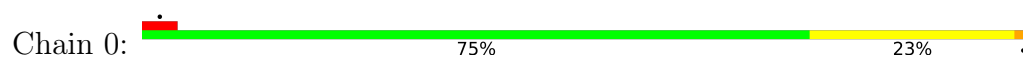


Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
67	8	1	9	6	2	1	0

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 L34



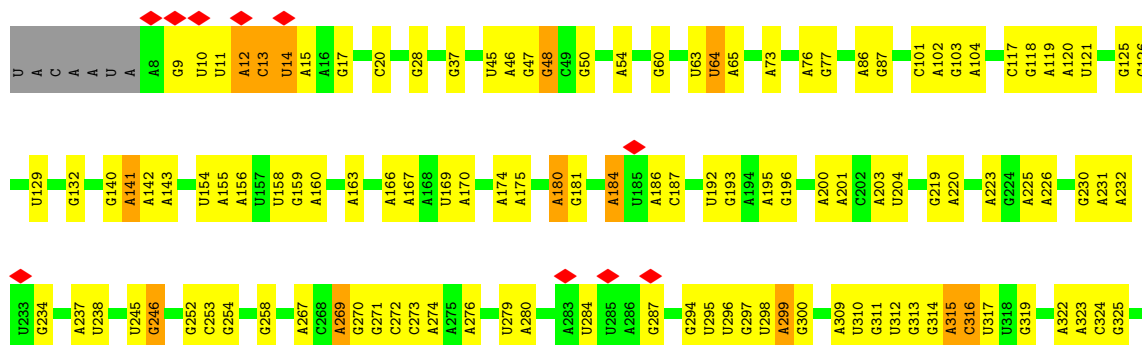
- Molecule 2: 50S ribosomal protein L35

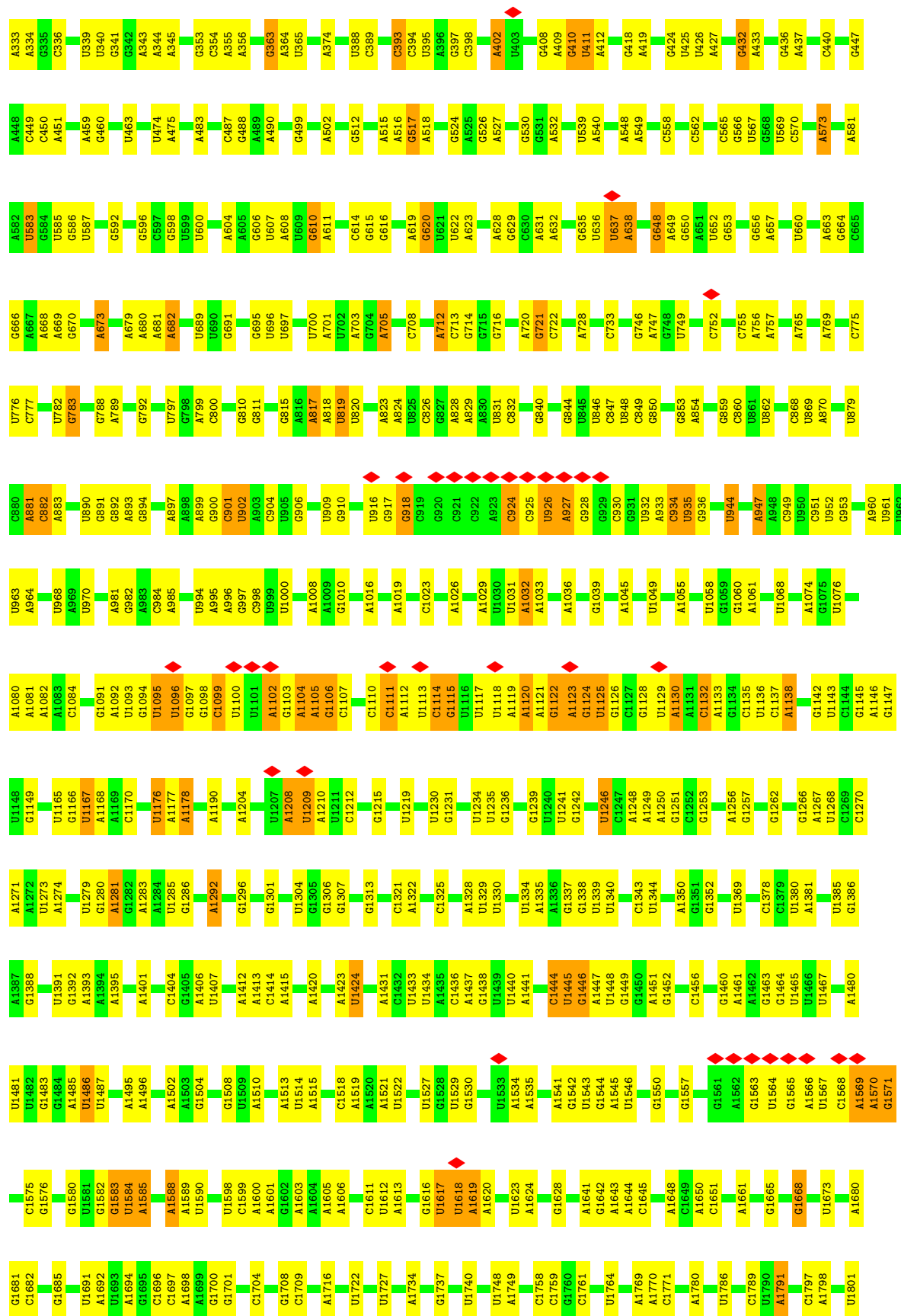


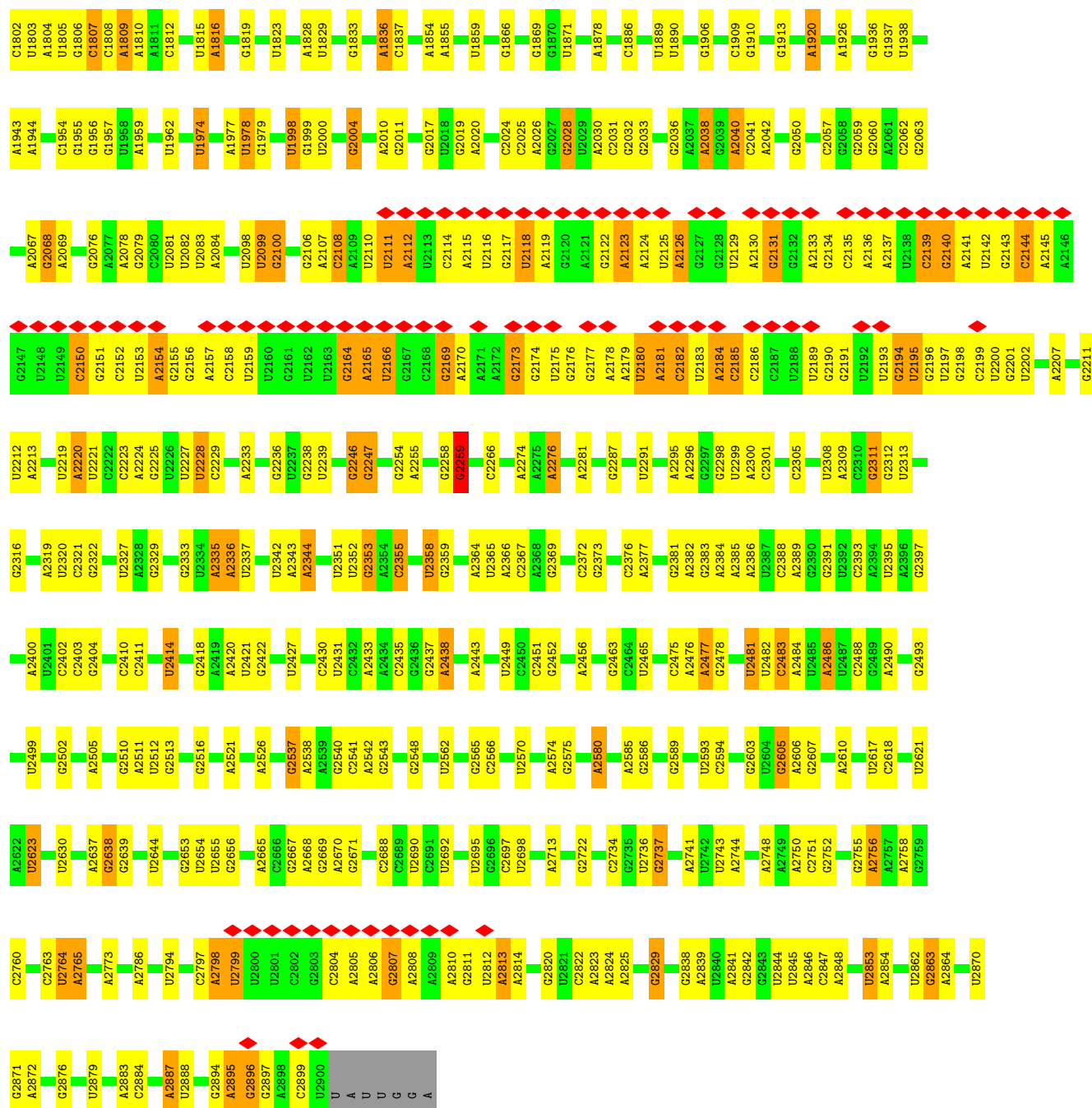
- Molecule 3: 50S ribosomal protein L36



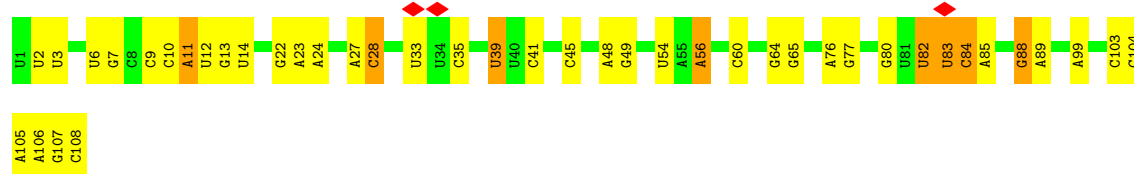
- Molecule 4: 23S ribosomal RNA

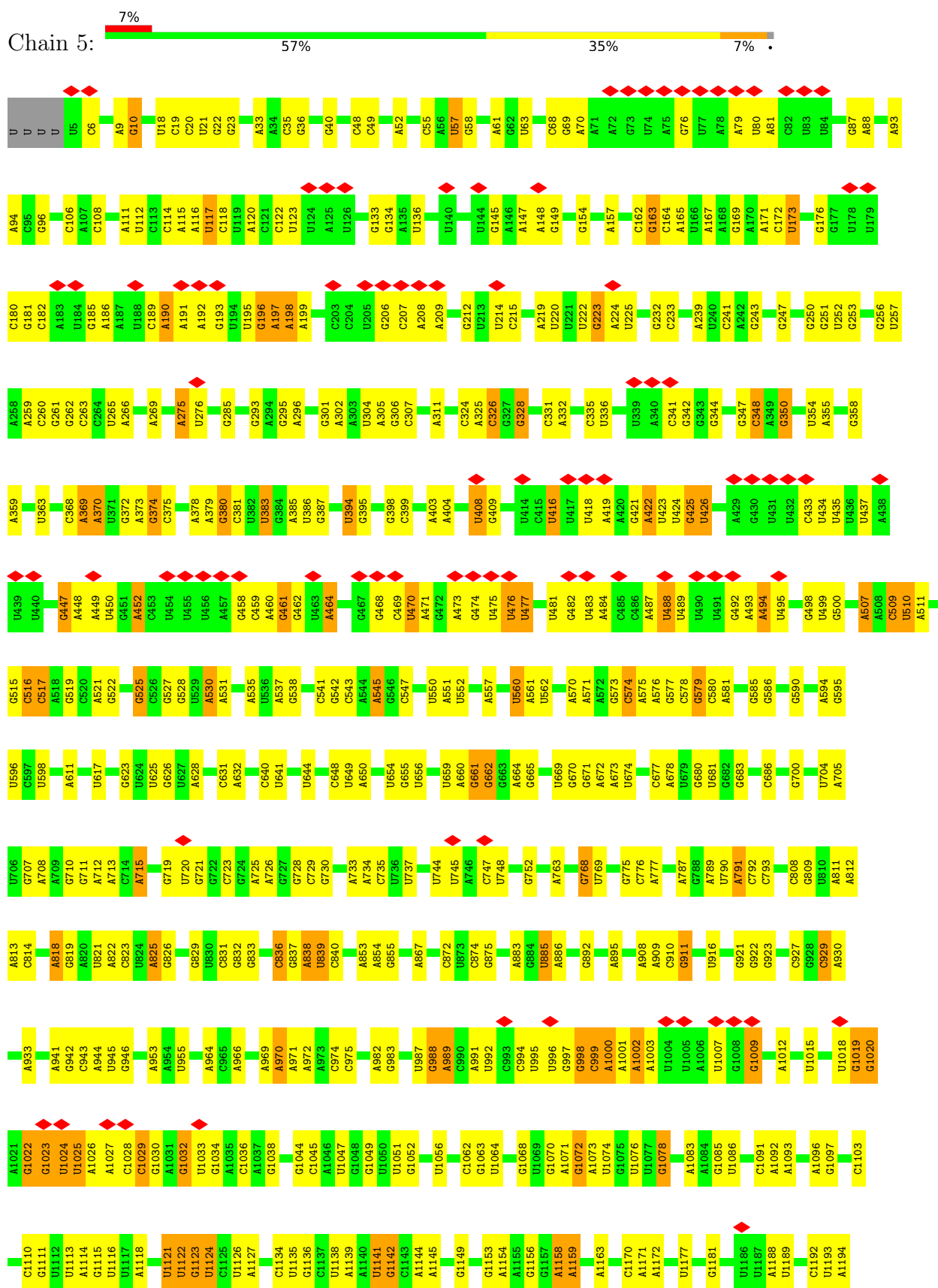


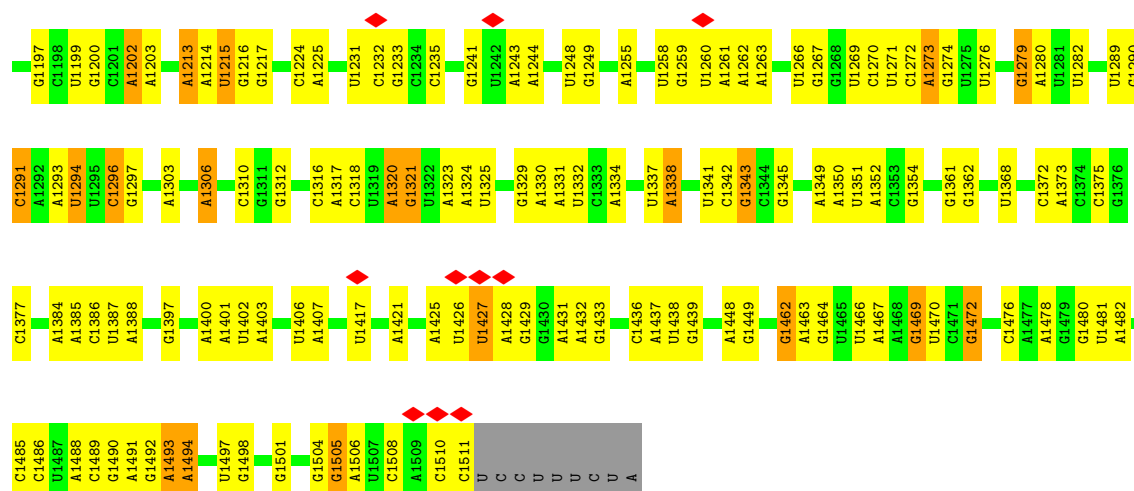




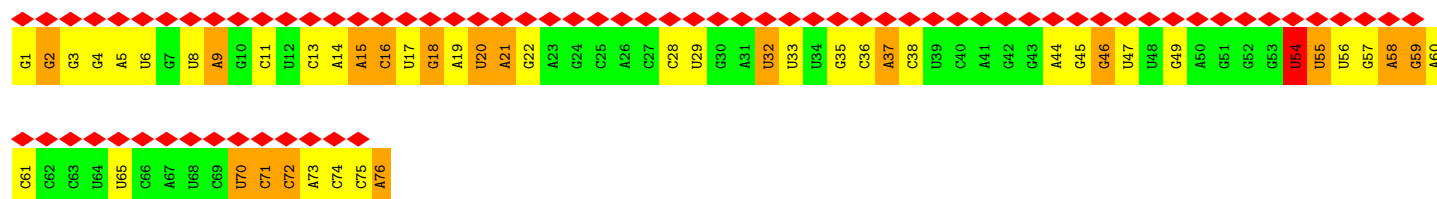
• Molecule 5: 5S ribosomal RNA



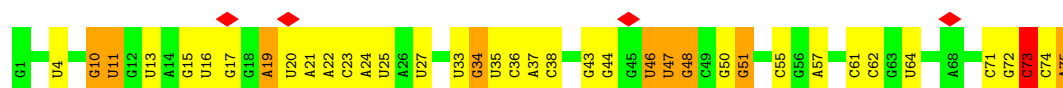




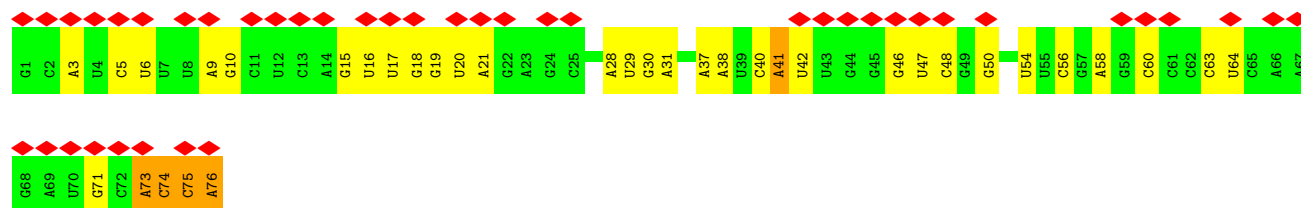
• Molecule 7: tRNA-Ala (E-site)



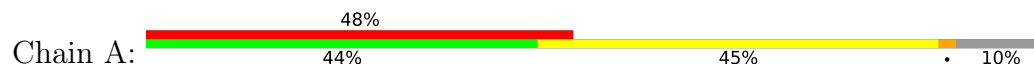
• Molecule 8: tRNA-Asp (P-site)

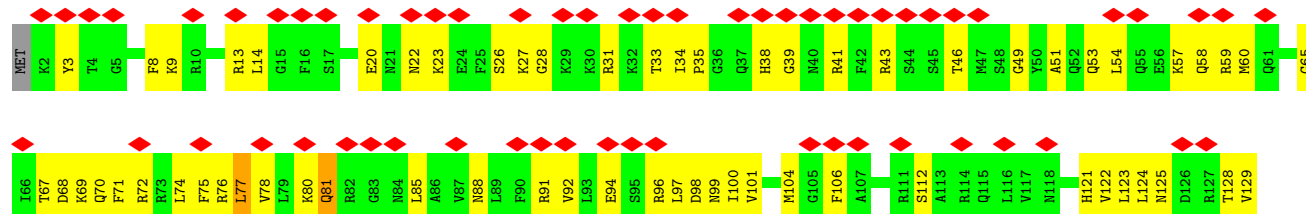


• Molecule 9: tRNA-Lys (A-site)



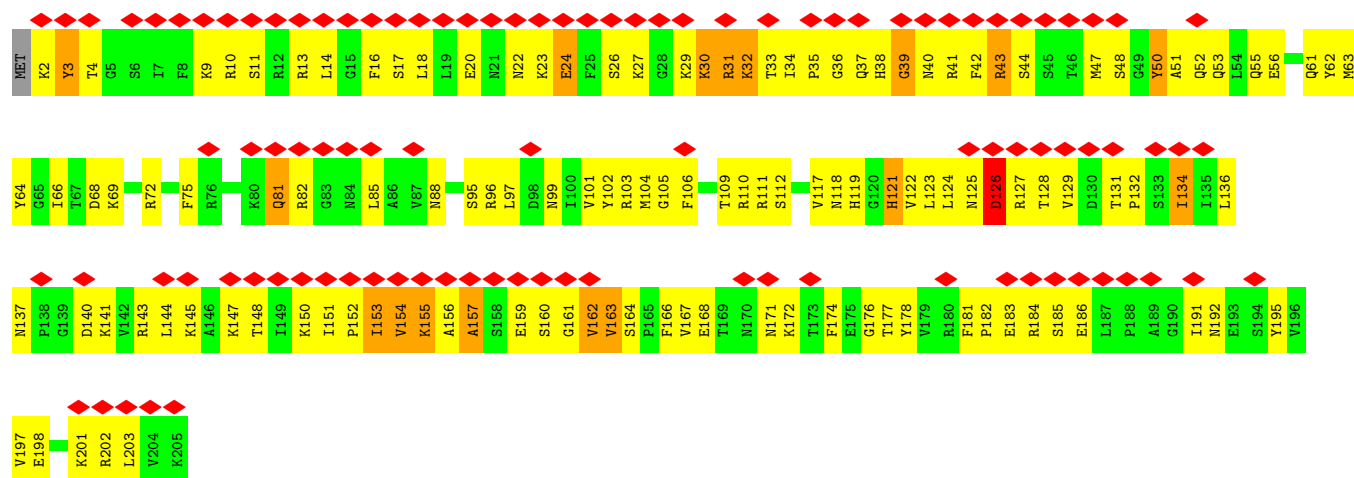
• Molecule 10: 30S ribosomal protein S2



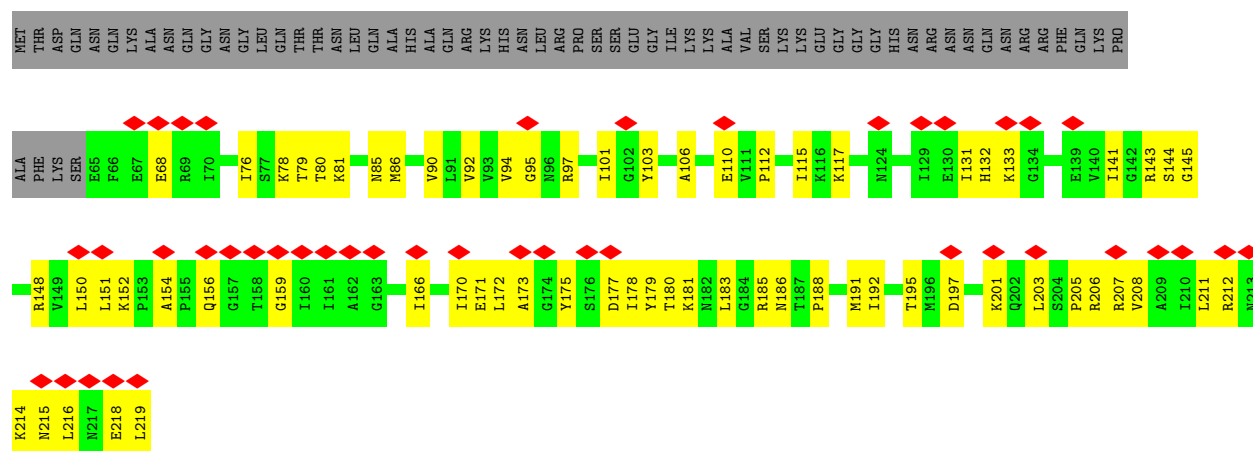




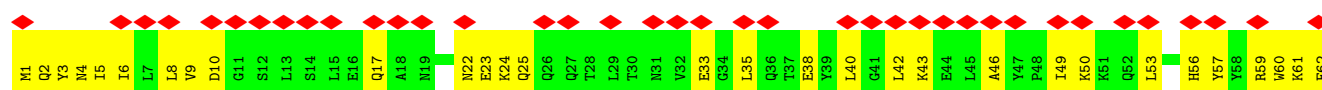
• Molecule 12: 30S ribosomal protein S4

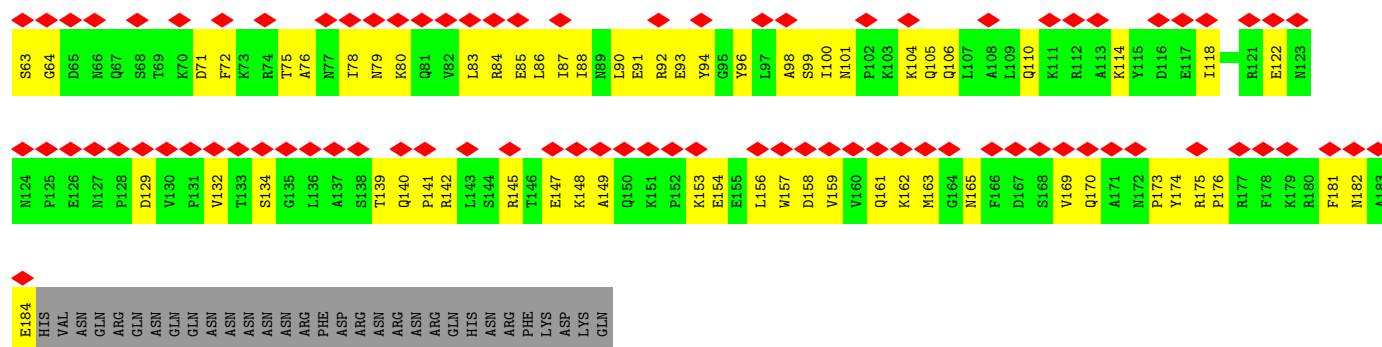


• Molecule 13: 30S ribosomal protein S5

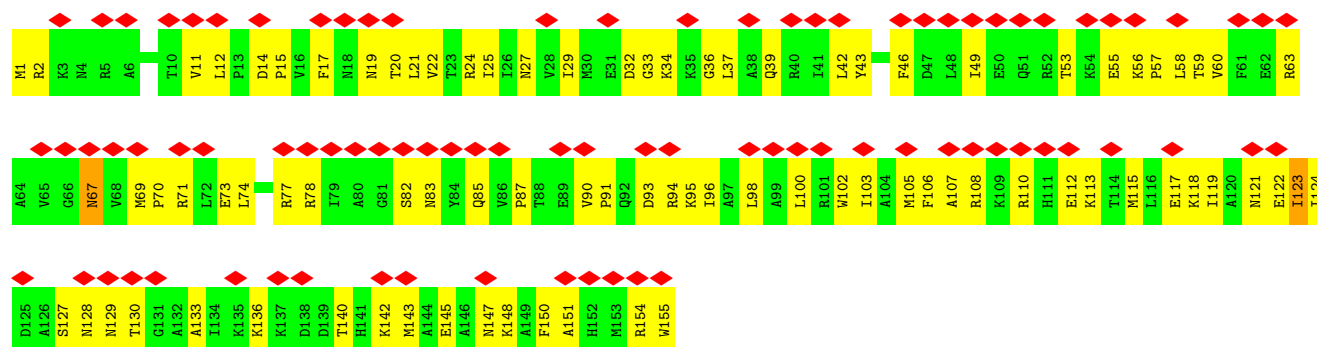


• Molecule 14: 30S ribosomal protein S6

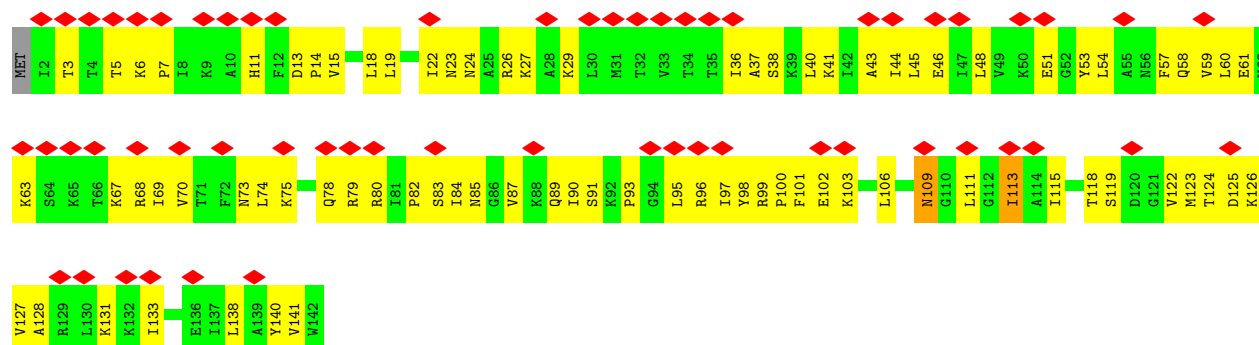
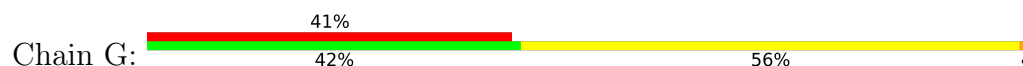




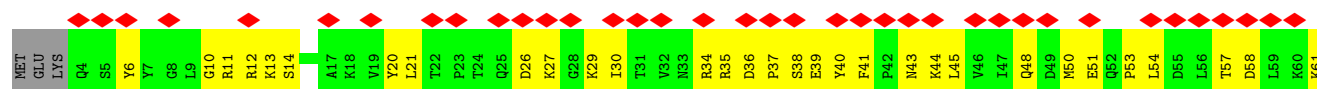
• Molecule 15: 30S ribosomal protein S7

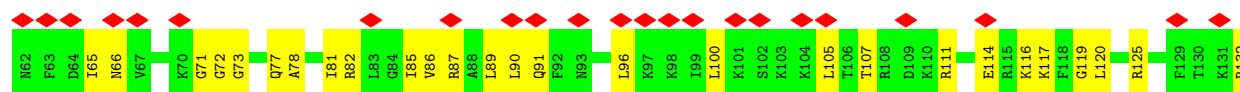


• Molecule 16: 30S ribosomal protein S8

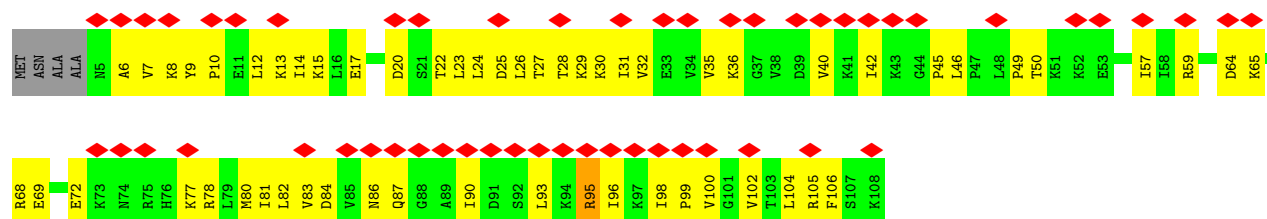
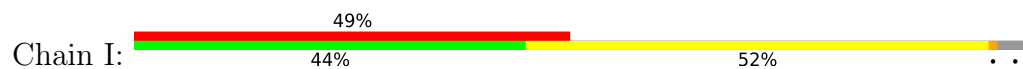


• Molecule 17: 30S ribosomal protein S9

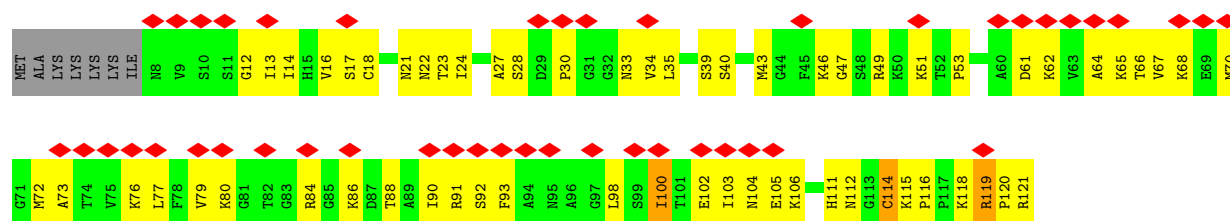
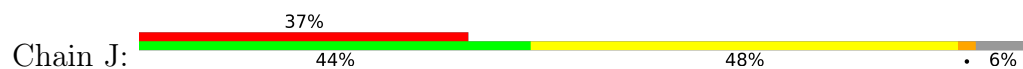




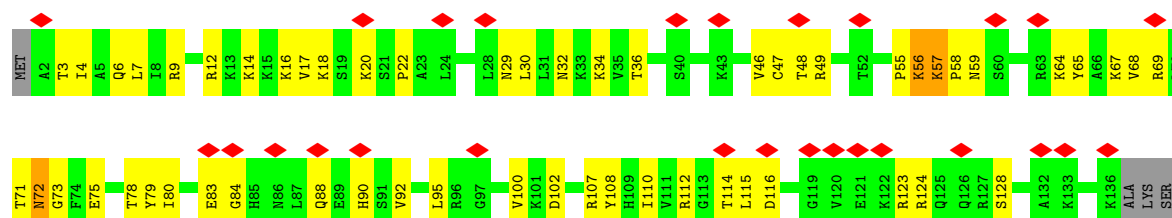
- Molecule 18: 30S ribosomal protein S10



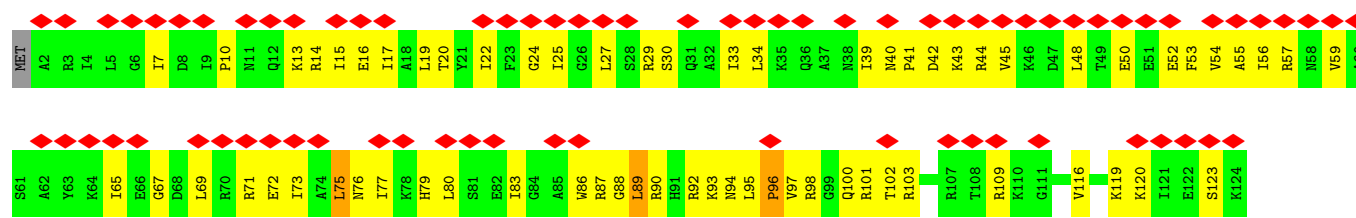
- Molecule 19: 30S ribosomal protein S11



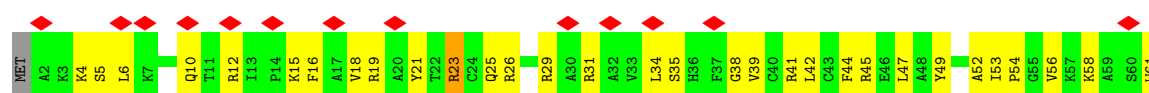
- Molecule 20: 30S ribosomal protein S12



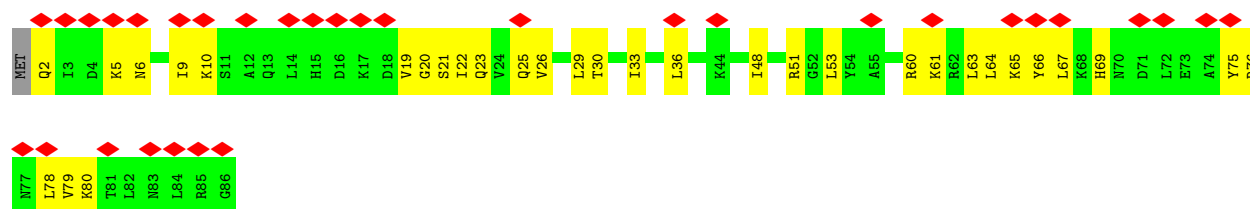
- Molecule 21: 30S ribosomal protein S13



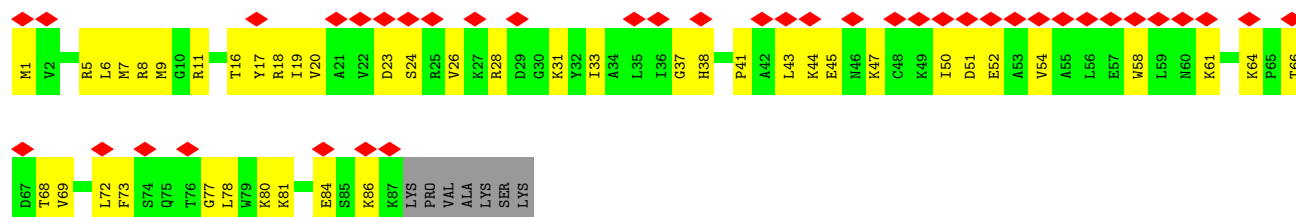
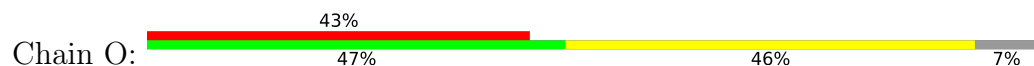
- Molecule 22: 30S ribosomal protein S14 type Z



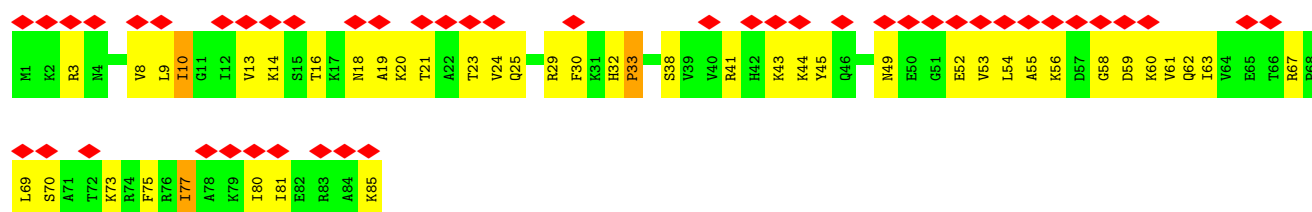
• Molecule 23: 30S ribosomal protein S15



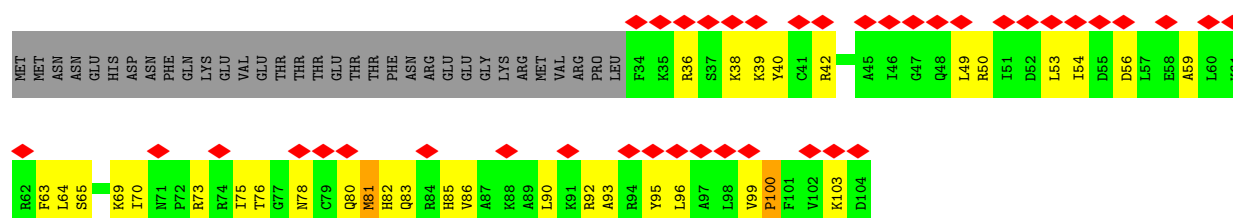
• Molecule 24: 30S ribosomal protein S16



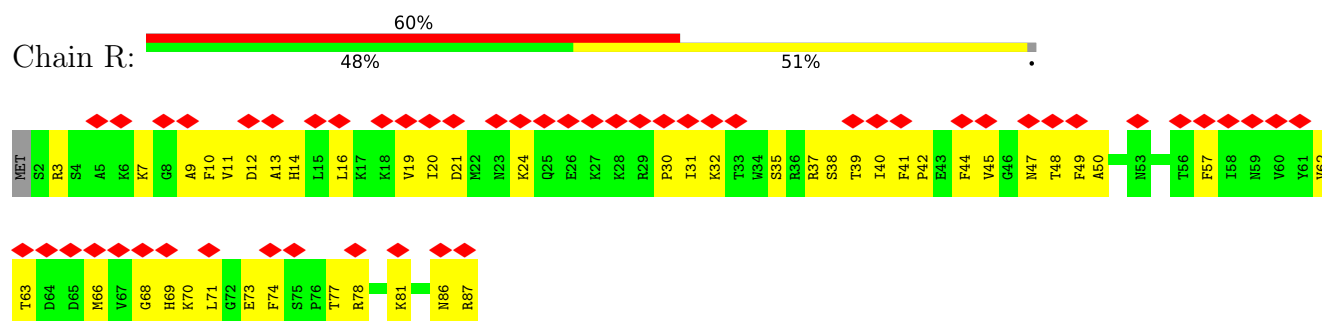
• Molecule 25: 30S ribosomal protein S17



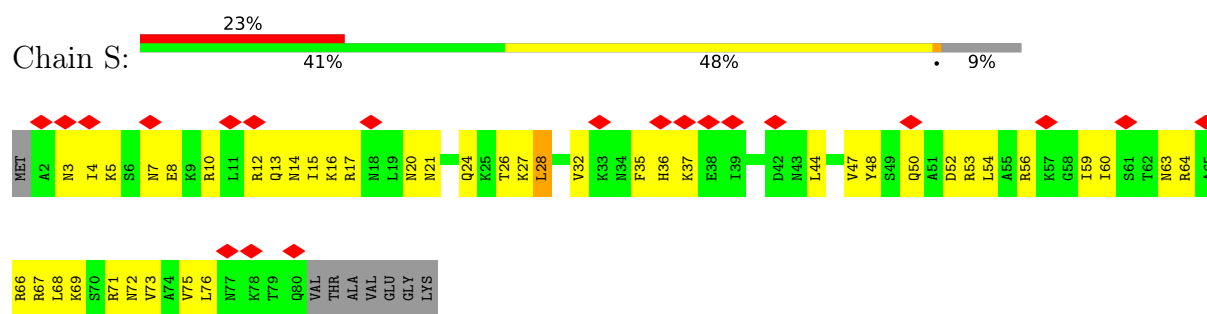
• Molecule 26: 30S ribosomal protein S18



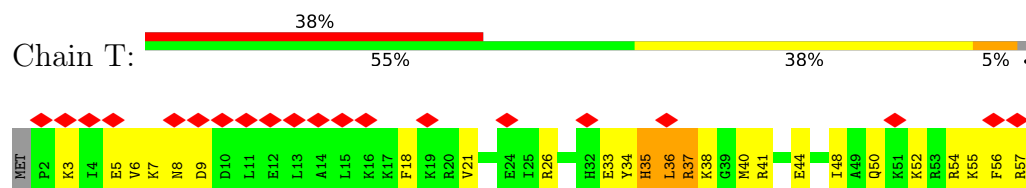
- Molecule 27: 30S ribosomal protein S19



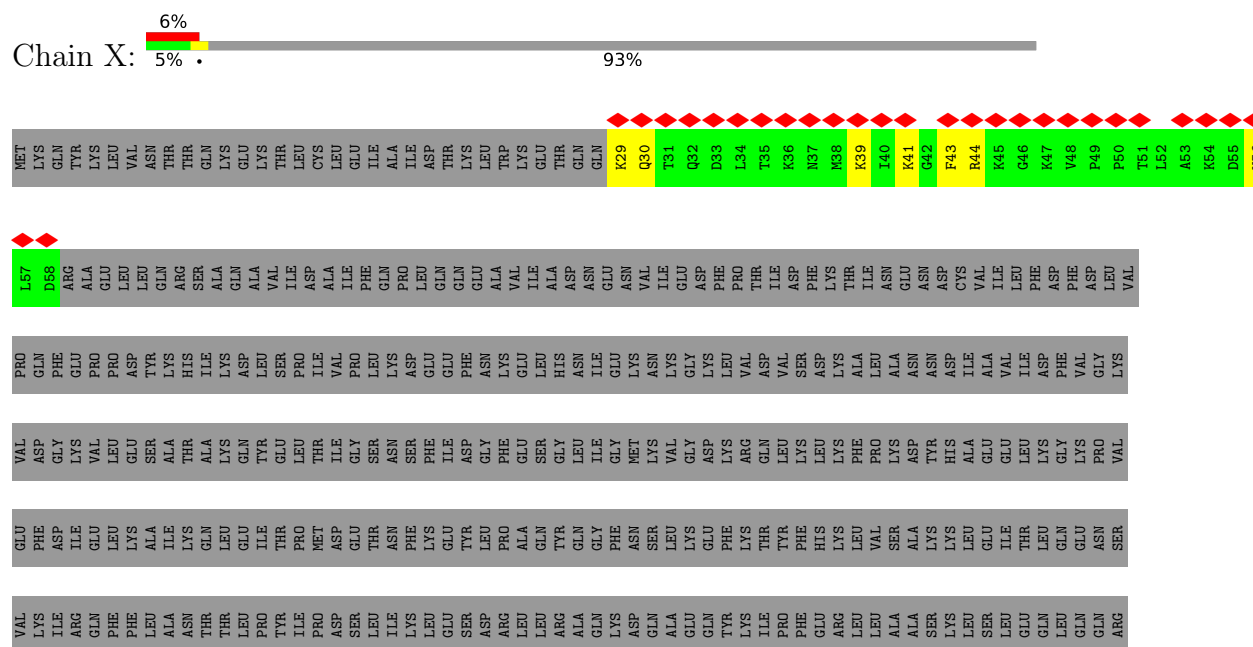
- Molecule 28: 30S ribosomal protein S20



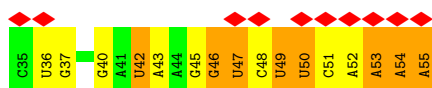
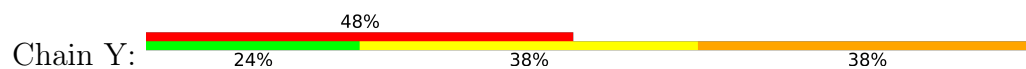
- Molecule 29: 30S ribosomal protein S21



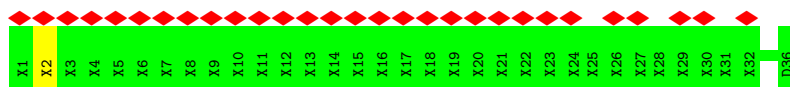
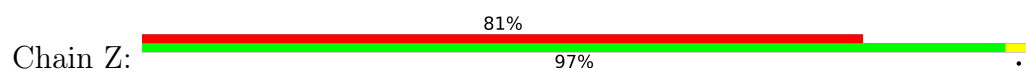
- Molecule 30: Trigger factor



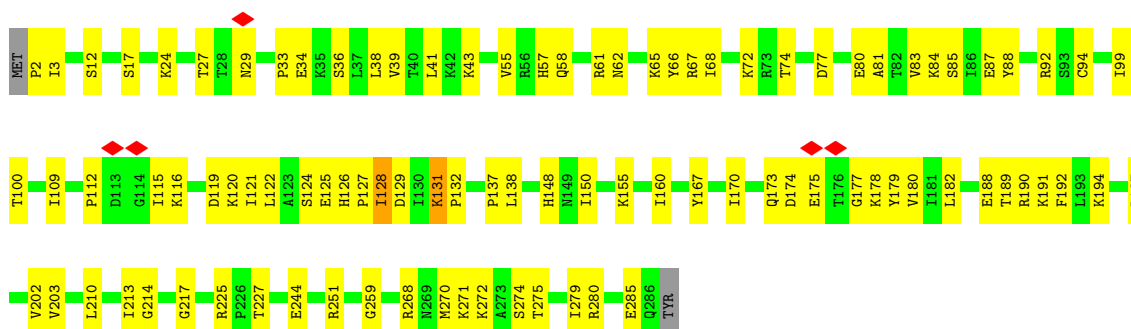
- Molecule 31: mRNA



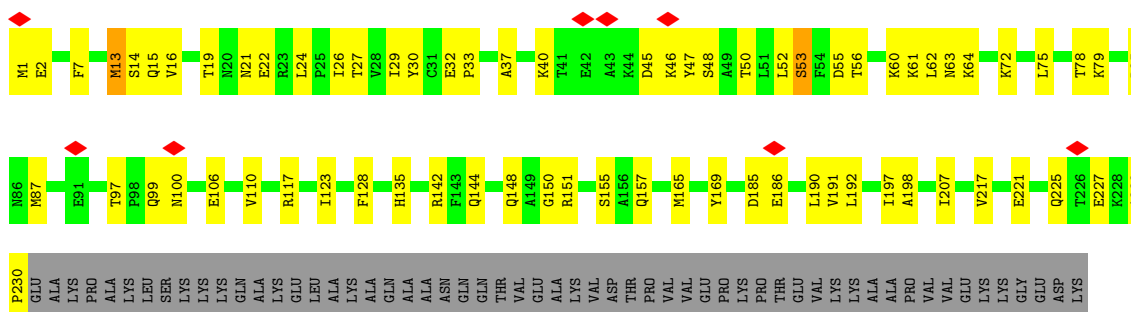
- Molecule 32: Nascent chain



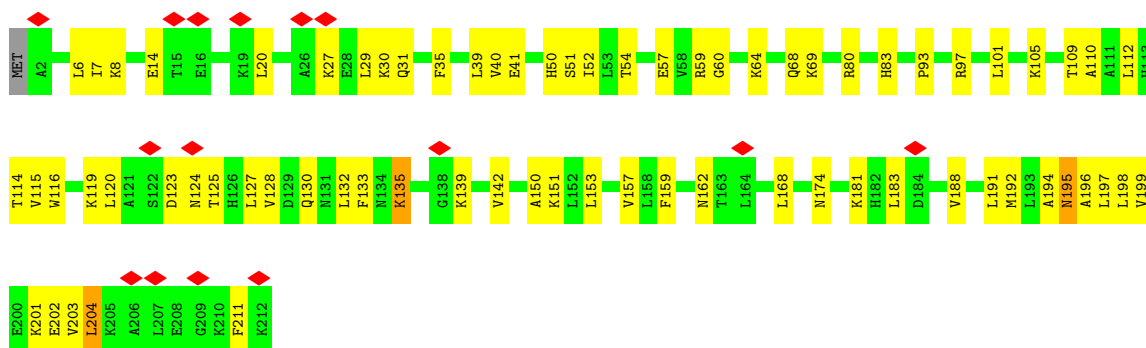
- Molecule 33: 50S ribosomal protein L2



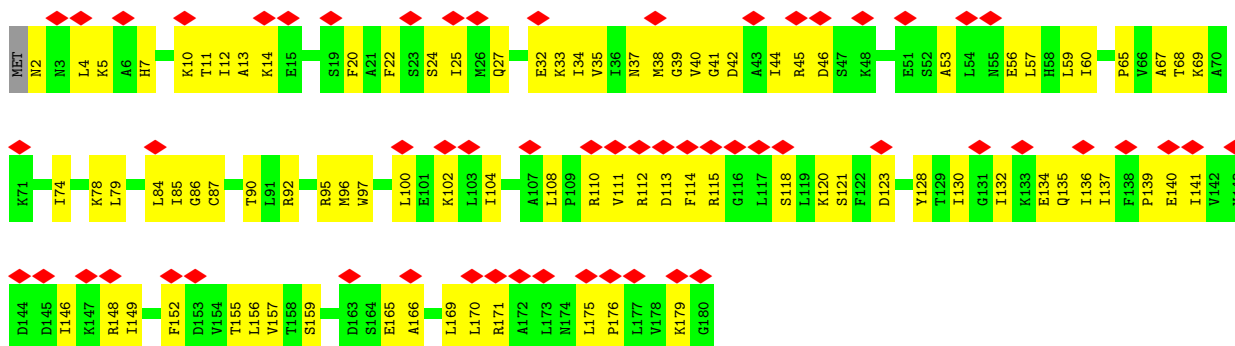
- Molecule 34: 50S ribosomal protein L3



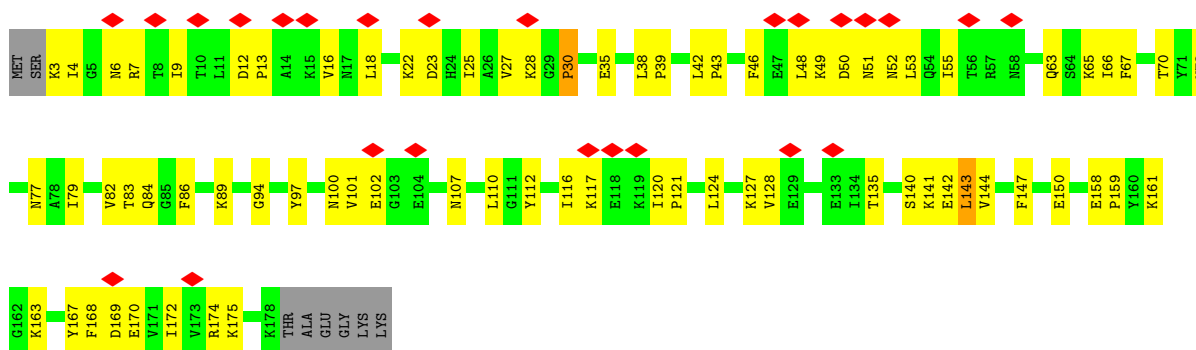
- Molecule 35: 50S ribosomal protein L4



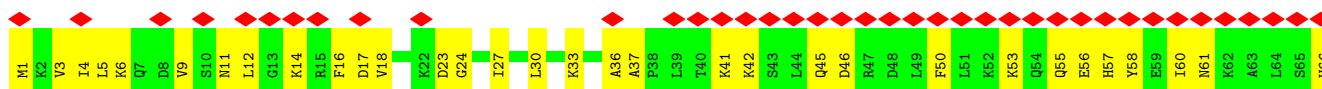
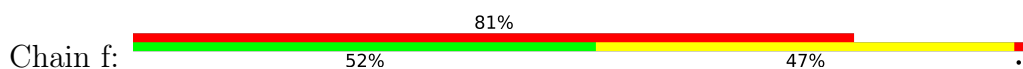
• Molecule 36: 50S ribosomal protein L5

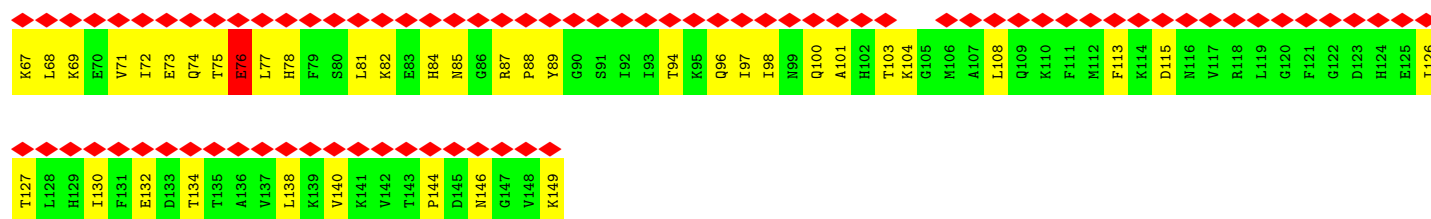


• Molecule 37: 50S ribosomal protein L6

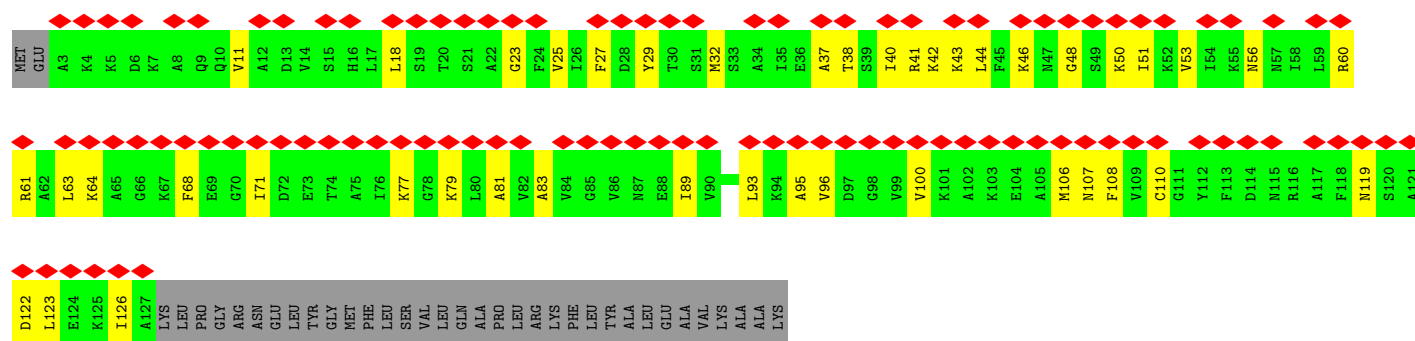


• Molecule 38: 50S ribosomal protein L9

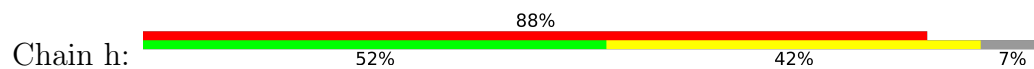




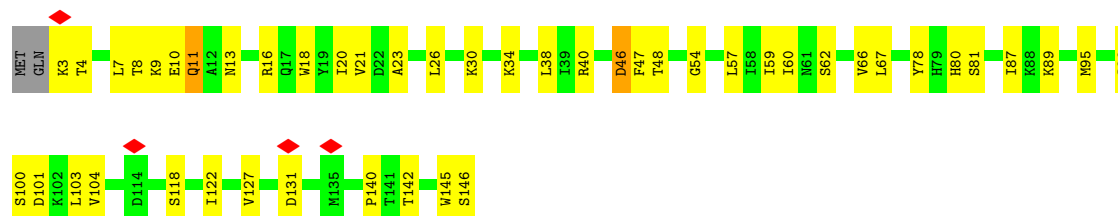
• Molecule 39: 50S ribosomal protein L10



• Molecule 40: 50S ribosomal protein L11

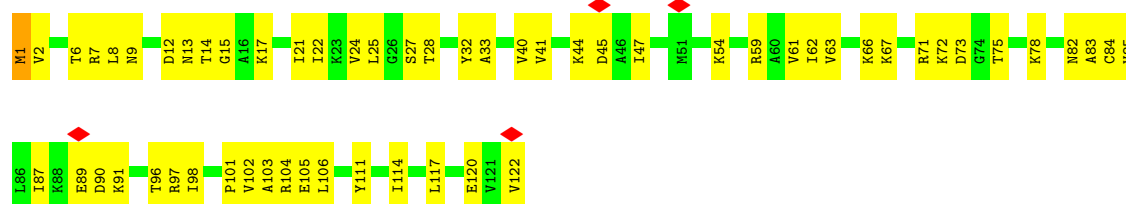


• Molecule 41: 50S ribosomal protein L13



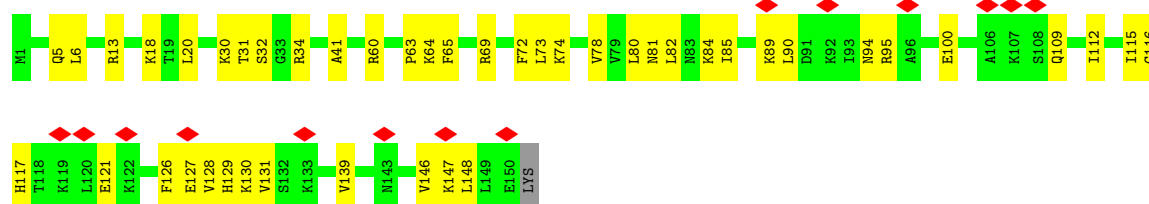
• Molecule 42: 50S ribosomal protein L14

Chain j: 



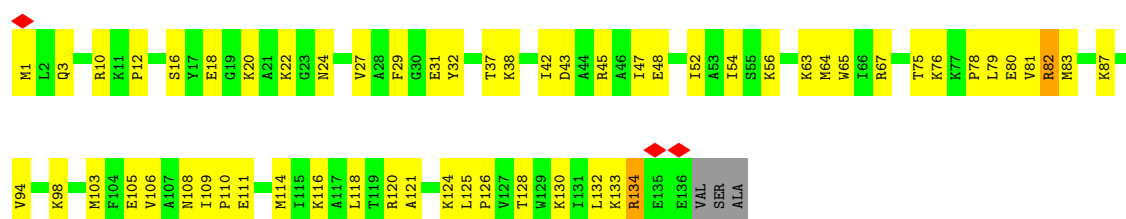
- Molecule 43: 50S ribosomal protein L15

Chain k: 



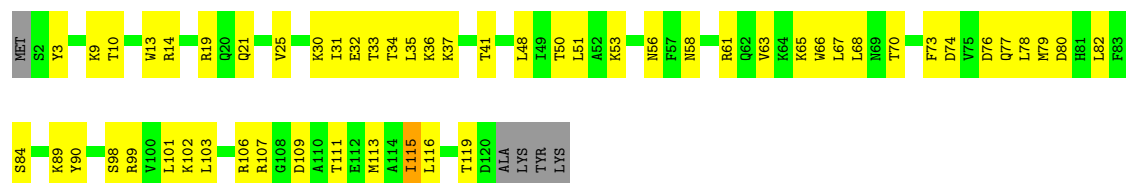
- Molecule 44: 50S ribosomal protein L16

Chain l: 



- Molecule 45: 50S ribosomal protein L17

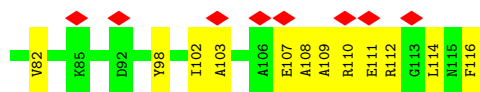
Chain m: 



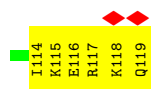
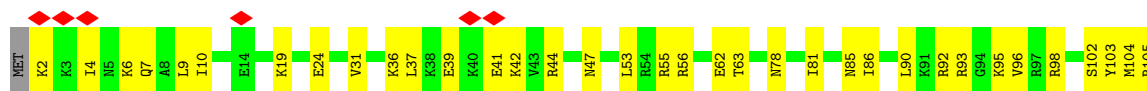
- Molecule 46: 50S ribosomal protein L18

Chain n: 





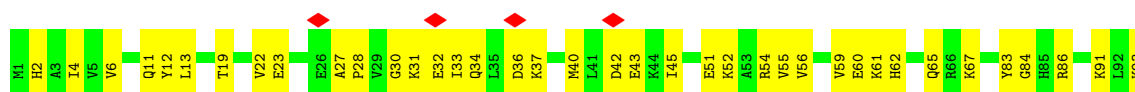
• Molecule 47: 50S ribosomal protein L19



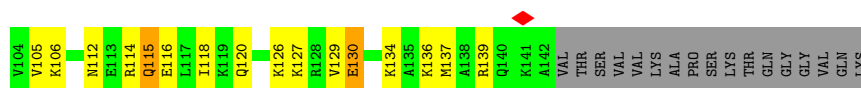
• Molecule 48: 50S ribosomal protein L20



• Molecule 49: 50S ribosomal protein L21

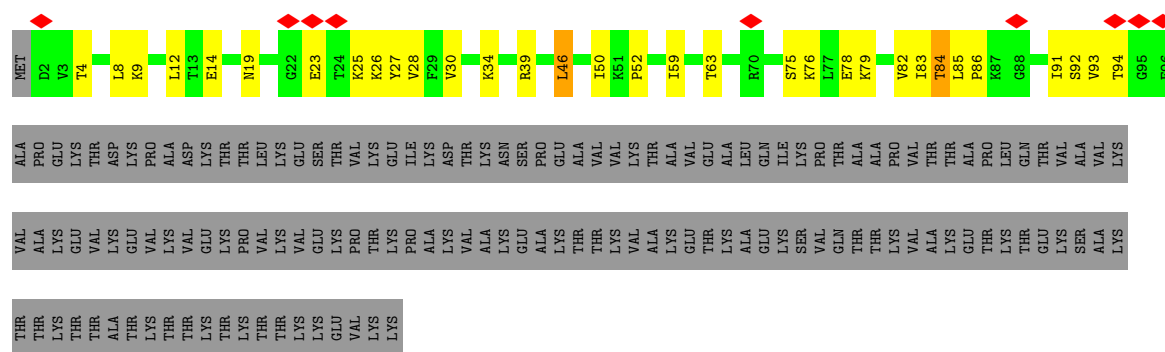


• Molecule 50: 50S ribosomal protein L22



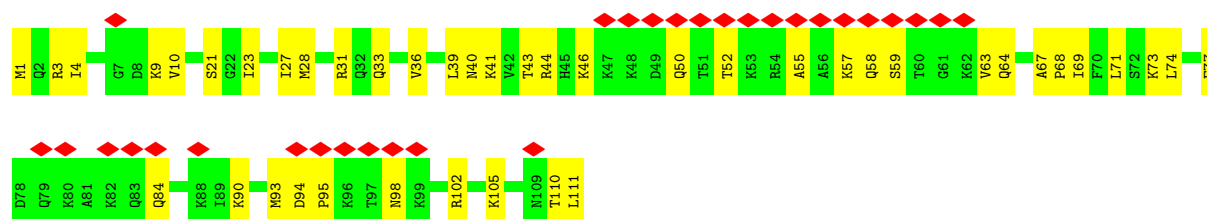
• Molecule 51: 50S ribosomal protein L23

Chain s: 



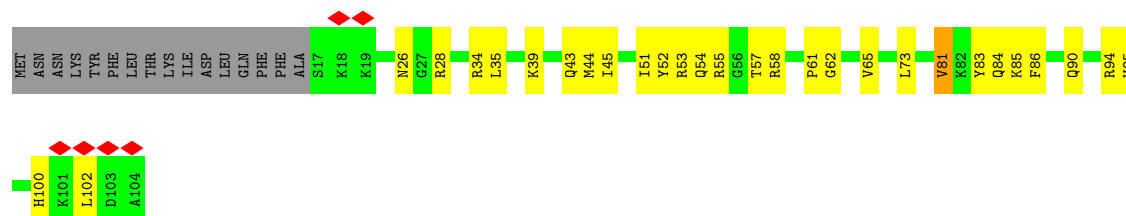
• Molecule 52: 50S ribosomal protein L24

Chain t: 



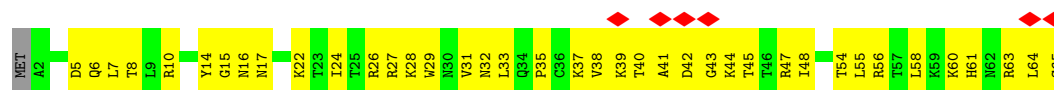
• Molecule 53: 50S ribosomal protein L27

Chain u: 



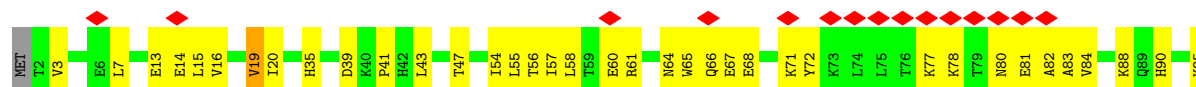
• Molecule 54: 50S ribosomal protein L28

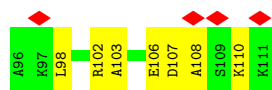
Chain v: 



• Molecule 55: 50S ribosomal protein L29

Chain w: 





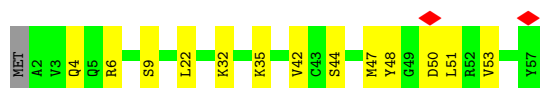
- Molecule 56: 50S ribosomal protein L31

Chain x: 73%
63% 35%



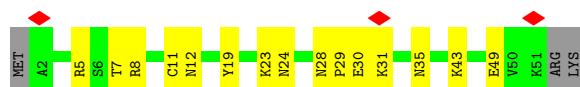
- Molecule 57: 50S ribosomal protein L32

Chain y: 75% 23%



- Molecule 58: 50S ribosomal protein L33 1

Chain z: 6% 66% 28% 6%



4 Experimental information

Property	Value	Source
EM reconstruction method	SUBTOMOGRAM AVERAGING	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of subtomograms used	10154	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; CTF estimation and 3D CTF correction are done in Warp	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	137	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3250	Depositor
Magnification	64000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	7.019	Depositor
Minimum map value	-3.485	Depositor
Average map value	0.033	Depositor
Map value standard deviation	0.273	Depositor
Recommended contour level	0.99	Depositor
Map size ($\text{\AA}$)	435.328, 435.328, 435.328	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.7005, 1.7005, 1.7005	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CLM, 1MG, MG, 7MG, SPM, 5MC, SPD, ZN, MA6, OMG, B8T, N2P, 2MA, K, PUT

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	0.19	0/395	0.47	0/518
2	1	0.17	0/484	0.37	0/637
3	2	0.19	0/306	0.53	0/401
4	3	0.09	0/69363	0.22	5/108161 (0.0%)
5	4	0.08	0/2578	0.19	0/4016
6	5	0.09	0/35992	0.21	0/56111
7	6	0.31	0/1810	0.54	2/2817 (0.1%)
8	7	0.19	0/1785	0.37	1/2779 (0.0%)
9	8	0.23	1/1804 (0.1%)	0.34	0/2807
10	A	0.27	0/2172	0.68	1/2934 (0.0%)
11	B	0.26	0/1863	0.67	3/2516 (0.1%)
12	C	0.20	0/1700	0.59	0/2278
12	U	0.41	1/1700 (0.1%)	0.90	8/2278 (0.4%)
13	D	0.21	0/1206	0.57	0/1616
14	E	0.31	0/1536	0.73	1/2072 (0.0%)
15	F	0.25	0/1274	0.68	0/1710
16	G	0.29	0/1126	0.86	6/1517 (0.4%)
17	H	0.33	0/1056	0.72	0/1409
18	I	0.23	0/843	0.65	3/1132 (0.3%)
19	J	0.29	0/844	0.67	2/1136 (0.2%)
20	K	0.22	0/1089	0.57	0/1461
21	L	0.24	0/1002	0.74	3/1340 (0.2%)
22	M	0.26	0/483	0.64	0/643
23	N	0.17	0/695	0.59	0/926
24	O	0.28	0/718	0.76	1/962 (0.1%)
25	P	0.20	0/702	0.58	1/934 (0.1%)
26	Q	0.36	0/601	0.86	1/801 (0.1%)
27	R	0.34	0/716	0.66	0/958
28	S	0.23	0/645	0.60	0/857
29	T	0.37	0/524	0.77	0/685
30	X	0.16	0/245	0.45	0/325
31	Y	0.17	0/498	0.37	0/773

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	Z	0.90	0/26	1.70	0/33
33	a	0.18	0/2267	0.50	1/3044 (0.0%)
34	b	0.21	0/1803	0.46	1/2424 (0.0%)
35	c	0.17	0/1681	0.46	0/2257
36	d	0.21	0/1437	0.57	0/1931
37	e	0.19	0/1420	0.52	1/1912 (0.1%)
38	f	0.32	0/1233	0.64	0/1653
39	g	0.16	0/960	0.55	0/1284
40	h	0.18	0/968	0.60	3/1298 (0.2%)
41	i	0.23	0/1186	0.57	1/1592 (0.1%)
42	j	0.23	0/953	0.55	0/1275
43	k	0.20	0/1187	0.52	0/1581
44	l	0.30	0/1104	0.67	1/1481 (0.1%)
45	m	0.17	0/973	0.46	0/1309
46	n	0.24	0/927	0.66	0/1239
47	o	0.20	0/976	0.57	0/1296
48	p	0.21	0/996	0.50	0/1325
49	q	0.18	0/828	0.53	0/1111
50	r	0.24	0/1100	0.63	3/1471 (0.2%)
51	s	0.17	0/752	0.46	0/1015
52	t	0.22	0/878	0.64	4/1165 (0.3%)
53	u	0.19	0/678	0.50	0/902
54	v	0.35	0/526	0.73	0/703
55	w	0.15	0/916	0.45	0/1222
56	x	0.18	0/776	0.58	0/1033
57	y	0.16	0/457	0.53	0/601
58	z	0.22	0/412	0.49	0/547
All	All	0.16	2/165175 (0.0%)	0.38	53/246214 (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
12	U	0	2
22	M	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	U	31	ARG	CA-C	9.23	1.68	1.52
9	8	76	A	C6-N6	6.01	1.46	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	A	173	PRO	CA-N-CD	-11.78	95.50	112.00
26	Q	100	PRO	CA-N-CD	-10.62	97.13	112.00
16	G	14	PRO	CA-N-CD	-8.47	100.14	112.00
21	L	96	PRO	CA-N-CD	-7.95	100.87	112.00
7	6	32	U	P-O3'-C3'	-7.88	108.38	120.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
22	M	23	ARG	Sidechain
12	U	121	HIS	Peptide
12	U	153	ILE	Peptide

5.2 Too-close contacts [i](#)

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	0	392	0	442	11	0
2	1	477	0	530	17	0
3	2	304	0	347	10	0
4	3	61995	0	31115	619	0
5	4	2305	0	1164	17	0
6	5	32258	0	16206	463	0
7	6	1620	0	818	28	0
8	7	1599	0	805	20	0
9	8	1615	0	816	15	0
10	A	2138	0	2204	119	0
11	B	1835	0	1909	110	0
12	C	1669	0	1729	90	0
12	U	1669	0	1729	159	0
13	D	1191	0	1284	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	E	1509	0	1520	102	0
15	F	1254	0	1320	96	0
16	G	1110	0	1226	96	0
17	H	1040	0	1107	71	0
18	I	832	0	918	51	0
19	J	829	0	855	69	0
20	K	1071	0	1165	56	0
21	L	991	0	1061	66	0
22	M	474	0	505	48	0
23	N	689	0	746	29	0
24	O	705	0	755	43	0
25	P	693	0	753	40	0
26	Q	590	0	626	38	0
27	R	700	0	709	43	0
28	S	643	0	694	42	0
29	T	519	0	578	42	0
30	X	242	0	263	7	0
31	Y	446	0	227	12	0
32	Z	187	0	68	1	0
33	a	2225	0	2301	80	0
34	b	1769	0	1815	65	0
35	c	1654	0	1744	66	0
36	d	1416	0	1500	83	0
37	e	1396	0	1481	59	0
38	f	1210	0	1259	53	0
39	g	951	0	1001	36	0
40	h	959	0	1039	42	0
41	i	1164	0	1192	42	0
42	j	944	0	1019	46	0
43	k	1170	0	1274	43	0
44	l	1079	0	1134	63	0
45	m	958	0	1011	41	0
46	n	918	0	979	50	0
47	o	966	0	1042	39	0
48	p	981	0	1062	52	0
49	q	811	0	858	39	0
50	r	1091	0	1178	40	0
51	s	740	0	819	26	0
52	t	871	0	972	37	0
53	u	670	0	704	31	0
54	v	520	0	565	49	0
55	w	906	0	981	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	x	761	0	766	31	0
57	y	452	0	471	11	0
58	z	408	0	436	14	0
59	2	1	0	0	0	0
59	M	1	0	0	0	0
59	Q	1	0	0	0	0
59	x	1	0	0	0	0
59	y	1	0	0	0	0
59	z	1	0	0	0	0
60	3	20	0	11	1	0
61	3	1	0	0	0	0
62	3	221	0	0	0	0
62	4	1	0	0	0	0
62	5	90	0	0	0	0
62	6	1	0	0	0	0
62	7	2	0	0	0	0
62	8	2	0	0	0	0
62	K	1	0	0	0	0
62	P	1	0	0	0	0
62	Y	2	0	0	0	0
62	b	2	0	0	0	0
62	i	1	0	0	0	0
62	y	2	0	0	0	0
63	3	42	0	84	2	0
63	5	6	0	12	0	0
64	3	42	0	78	7	0
64	b	14	0	26	3	0
65	3	160	0	304	3	0
65	5	20	0	38	7	0
66	3	21	0	42	0	0
66	5	7	0	14	0	0
67	8	9	0	12	0	0
All	All	153255	0	105418	3310	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 13.

The worst 5 of 3310 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)
12:U:14:LEU:HA	54:v:63:ARG:NH1	1.67	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:a:131:LYS:HE3	33:a:132:PRO:HD3	1.35	1.07
6:5:974:C:H2'	22:M:19:ARG:HH12	1.18	1.04
47:o:37:LEU:HD11	47:o:44:ARG:HB3	1.42	1.01
6:5:172:C:H5''	28:S:66:ARG:HH11	1.21	1.00

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	0	46/48 (96%)	45 (98%)	1 (2%)	0	100	100
2	1	57/59 (97%)	57 (100%)	0	0	100	100
3	2	35/37 (95%)	35 (100%)	0	0	100	100
10	A	262/294 (89%)	236 (90%)	24 (9%)	2 (1%)	16	48
11	B	230/273 (84%)	203 (88%)	23 (10%)	4 (2%)	7	34
12	C	202/205 (98%)	167 (83%)	33 (16%)	2 (1%)	12	43
12	U	202/205 (98%)	147 (73%)	39 (19%)	16 (8%)	1	10
13	D	153/219 (70%)	143 (94%)	10 (6%)	0	100	100
14	E	182/215 (85%)	156 (86%)	25 (14%)	1 (0%)	24	56
15	F	153/155 (99%)	139 (91%)	12 (8%)	2 (1%)	9	38
16	G	139/142 (98%)	133 (96%)	6 (4%)	0	100	100
17	H	127/132 (96%)	113 (89%)	14 (11%)	0	100	100
18	I	102/108 (94%)	89 (87%)	12 (12%)	1 (1%)	12	43
19	J	112/121 (93%)	106 (95%)	5 (4%)	1 (1%)	14	45
20	K	133/139 (96%)	116 (87%)	14 (10%)	3 (2%)	5	30
21	L	121/124 (98%)	109 (90%)	10 (8%)	2 (2%)	7	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
22	M	58/61 (95%)	56 (97%)	2 (3%)	0	100	100
23	N	83/86 (96%)	78 (94%)	5 (6%)	0	100	100
24	O	85/94 (90%)	79 (93%)	6 (7%)	0	100	100
25	P	83/85 (98%)	73 (88%)	9 (11%)	1 (1%)	10	40
26	Q	69/104 (66%)	64 (93%)	5 (7%)	0	100	100
27	R	84/87 (97%)	78 (93%)	6 (7%)	0	100	100
28	S	77/87 (88%)	74 (96%)	2 (3%)	1 (1%)	9	38
29	T	57/60 (95%)	54 (95%)	3 (5%)	0	100	100
30	X	28/444 (6%)	25 (89%)	3 (11%)	0	100	100
32	Z	3/36 (8%)	3 (100%)	0	0	100	100
33	a	283/287 (99%)	269 (95%)	13 (5%)	1 (0%)	30	60
34	b	228/287 (79%)	219 (96%)	9 (4%)	0	100	100
35	c	209/212 (99%)	196 (94%)	13 (6%)	0	100	100
36	d	177/180 (98%)	171 (97%)	6 (3%)	0	100	100
37	e	174/184 (95%)	161 (92%)	13 (8%)	0	100	100
38	f	147/149 (99%)	133 (90%)	12 (8%)	2 (1%)	9	37
39	g	123/161 (76%)	119 (97%)	4 (3%)	0	100	100
40	h	126/137 (92%)	121 (96%)	5 (4%)	0	100	100
41	i	142/146 (97%)	134 (94%)	8 (6%)	0	100	100
42	j	120/122 (98%)	115 (96%)	5 (4%)	0	100	100
43	k	148/151 (98%)	138 (93%)	10 (7%)	0	100	100
44	l	134/139 (96%)	132 (98%)	2 (2%)	0	100	100
45	m	117/124 (94%)	113 (97%)	4 (3%)	0	100	100
46	n	114/116 (98%)	108 (95%)	6 (5%)	0	100	100
47	o	116/119 (98%)	108 (93%)	8 (7%)	0	100	100
48	p	116/127 (91%)	114 (98%)	2 (2%)	0	100	100
49	q	97/100 (97%)	92 (95%)	5 (5%)	0	100	100
50	r	140/159 (88%)	136 (97%)	4 (3%)	0	100	100
51	s	93/237 (39%)	89 (96%)	4 (4%)	0	100	100
52	t	109/111 (98%)	103 (94%)	6 (6%)	0	100	100
53	u	86/104 (83%)	83 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
54	v	62/65 (95%)	57 (92%)	5 (8%)	0	100	100
55	w	108/111 (97%)	104 (96%)	4 (4%)	0	100	100
56	x	94/97 (97%)	70 (74%)	22 (23%)	2 (2%)	5	31
57	y	54/57 (95%)	52 (96%)	2 (4%)	0	100	100
58	z	48/53 (91%)	47 (98%)	1 (2%)	0	100	100
All	All	6248/7355 (85%)	5762 (92%)	445 (7%)	41 (1%)	20	50

5 of 41 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
11	B	229	GLN
11	B	230	PRO
20	K	56	LYS
25	P	54	LEU
12	U	39	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	41/41 (100%)	40 (98%)	1 (2%)	43	61
2	1	51/51 (100%)	51 (100%)	0	100	100
3	2	35/35 (100%)	34 (97%)	1 (3%)	37	57
10	A	238/262 (91%)	235 (99%)	3 (1%)	61	71
11	B	195/232 (84%)	195 (100%)	0	100	100
12	C	182/183 (100%)	181 (100%)	1 (0%)	81	80
12	U	182/183 (100%)	179 (98%)	3 (2%)	55	68
13	D	125/178 (70%)	125 (100%)	0	100	100
14	E	165/196 (84%)	163 (99%)	2 (1%)	63	72
15	F	132/132 (100%)	130 (98%)	2 (2%)	57	68
16	G	123/124 (99%)	121 (98%)	2 (2%)	55	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	H	112/115 (97%)	112 (100%)	0	100	100
18	I	97/99 (98%)	97 (100%)	0	100	100
19	J	91/97 (94%)	89 (98%)	2 (2%)	45	63
20	K	117/120 (98%)	116 (99%)	1 (1%)	70	75
21	L	104/105 (99%)	103 (99%)	1 (1%)	68	74
22	M	47/48 (98%)	47 (100%)	0	100	100
23	N	77/78 (99%)	77 (100%)	0	100	100
24	O	76/82 (93%)	76 (100%)	0	100	100
25	P	75/75 (100%)	73 (97%)	2 (3%)	39	59
26	Q	62/94 (66%)	61 (98%)	1 (2%)	55	68
27	R	76/77 (99%)	76 (100%)	0	100	100
28	S	71/77 (92%)	69 (97%)	2 (3%)	38	58
29	T	55/56 (98%)	52 (94%)	3 (6%)	19	46
30	X	27/406 (7%)	27 (100%)	0	100	100
32	Z	2/2 (100%)	2 (100%)	0	100	100
33	a	241/243 (99%)	241 (100%)	0	100	100
34	b	187/233 (80%)	184 (98%)	3 (2%)	55	68
35	c	183/184 (100%)	179 (98%)	4 (2%)	45	63
36	d	153/154 (99%)	153 (100%)	0	100	100
37	e	153/159 (96%)	151 (99%)	2 (1%)	61	71
38	f	134/134 (100%)	133 (99%)	1 (1%)	76	77
39	g	100/129 (78%)	100 (100%)	0	100	100
40	h	102/110 (93%)	102 (100%)	0	100	100
41	i	126/128 (98%)	124 (98%)	2 (2%)	55	68
42	j	103/103 (100%)	101 (98%)	2 (2%)	50	65
43	k	125/126 (99%)	124 (99%)	1 (1%)	73	76
44	l	113/115 (98%)	111 (98%)	2 (2%)	51	67
45	m	105/109 (96%)	103 (98%)	2 (2%)	50	65
46	n	99/99 (100%)	98 (99%)	1 (1%)	68	74
47	o	104/105 (99%)	103 (99%)	1 (1%)	68	74
48	p	104/108 (96%)	104 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
49	q	90/91 (99%)	90 (100%)	0	100	100
50	r	118/132 (89%)	116 (98%)	2 (2%)	53	67
51	s	84/208 (40%)	81 (96%)	3 (4%)	31	54
52	t	96/96 (100%)	96 (100%)	0	100	100
53	u	70/85 (82%)	69 (99%)	1 (1%)	59	70
54	v	59/60 (98%)	59 (100%)	0	100	100
55	w	97/98 (99%)	96 (99%)	1 (1%)	68	74
56	x	85/86 (99%)	85 (100%)	0	100	100
57	y	48/49 (98%)	48 (100%)	0	100	100
58	z	47/50 (94%)	46 (98%)	1 (2%)	47	64
All	All	5484/6342 (86%)	5428 (99%)	56 (1%)	65	74

5 of 56 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
34	b	53	SER
58	z	43	LYS
37	e	143	LEU
55	w	19	VAL
50	r	115	GLN

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

Mol	Chain	Res	Type
33	a	57	HIS
53	u	26	ASN
35	c	76	GLN
50	r	112	ASN
54	v	16	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
31	Y	20/21 (95%)	14 (70%)	4 (20%)
4	3	2892/2907 (99%)	520 (17%)	21 (0%)
5	4	107/108 (99%)	29 (27%)	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
6	5	1506/1520 (99%)	254 (16%)	7 (0%)
7	6	75/76 (98%)	30 (40%)	6 (8%)
8	7	74/75 (98%)	20 (27%)	2 (2%)
9	8	75/76 (98%)	23 (30%)	0
All	All	4749/4783 (99%)	890 (18%)	40 (0%)

5 of 890 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	3	11	U
4	3	12	A
4	3	13	C
4	3	14	U
4	3	28	G

5 of 40 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
7	6	17	U
8	7	46	U
7	6	18	G
7	6	58	A
31	Y	49	U

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

8 non-standard protein/DNA/RNA residues are modelled in this entry.

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$
4	OMG	3	2259	4,62,8	23,26,27	2.63	6 (26%)	32,38,41	2.46	11 (34%)
6	7MG	5	525	6	23,26,27	3.62	11 (47%)	27,39,42	2.17	9 (33%)
6	B8T	5	1377	6	19,22,23	3.33	8 (42%)	25,31,34	0.84	1 (4%)
4	2MA	3	2511	4,62	22,25,26	4.13	9 (40%)	32,37,40	3.56	12 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MA6	5	1493	6	23,26,27	1.40	3 (13%)	33,38,41	3.24	14 (42%)
6	5MC	5	1375	6	19,22,23	4.56	8 (42%)	26,32,35	1.01	2 (7%)
6	MA6	5	1494	6	23,26,27	1.40	3 (13%)	33,38,41	3.26	14 (42%)
4	1MG	3	783	4	23,26,27	3.04	6 (26%)	33,39,42	2.32	8 (24%)

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
4	OMG	3	2259	4,62,8	-	3/9/27/28	0/3/3/3
6	7MG	5	525	6	-	2/7/37/38	0/3/3/3
6	B8T	5	1377	6	-	1/7/27/28	0/2/2/2
4	2MA	3	2511	4,62	-	1/7/25/26	0/3/3/3
6	MA6	5	1493	6	-	0/11/29/30	0/3/3/3
6	5MC	5	1375	6	-	0/7/25/26	0/2/2/2
6	MA6	5	1494	6	-	3/11/29/30	0/3/3/3
4	1MG	3	783	4	-	0/7/25/26	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	3	2511	2MA	C4-N3	11.99	1.49	1.34
6	5	1375	5MC	C6-C5	10.10	1.51	1.34
6	5	1375	5MC	C5-C4	9.23	1.51	1.44
4	3	783	1MG	C2-N3	8.63	1.47	1.33
6	5	525	7MG	C8-N9	8.60	1.51	1.45

The worst 5 of 71 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	3	2511	2MA	C1'-N9-C8	-12.59	99.15	127.09
4	3	2511	2MA	C4-N9-C1'	11.03	152.43	126.63
6	5	1494	MA6	N1-C6-N6	-10.42	104.16	116.86
6	5	1493	MA6	N1-C6-N6	-10.37	104.22	116.86
6	5	1493	MA6	C5-C6-N6	6.89	136.24	125.33

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	3	2259	OMG	O4'-C4'-C5'-O5'
4	3	2259	OMG	C3'-C4'-C5'-O5'
4	3	2259	OMG	C1'-C2'-O2'-CM2
6	5	525	7MG	O4'-C4'-C5'-O5'
6	5	1494	MA6	O4'-C4'-C5'-O5'

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	3	2259	OMG	2	0
6	5	1493	MA6	1	0
4	3	783	1MG	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 369 ligands modelled in this entry, 333 are monoatomic - leaving 36 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	$\# Z > 2$
63	PUT	3	3225	-	5,5,5	0.23	0	4,4,4	0.52	0
64	SPM	3	3243	-	13,13,13	0.34	0	12,12,12	0.91	0
67	LYS	8	103	9	7,8,9	0.87	0	3,8,10	0.32	0
63	PUT	3	3228	-	5,5,5	0.24	0	4,4,4	0.52	0
64	SPM	3	3229	-	13,13,13	0.36	0	12,12,12	0.90	0
65	SPD	3	3230	-	9,9,9	0.33	0	8,8,8	0.86	0
63	PUT	3	3244	-	5,5,5	0.24	0	4,4,4	0.52	0
65	SPD	5	1603	-	9,9,9	0.32	0	8,8,8	0.74	0
63	PUT	3	3224	-	5,5,5	0.24	0	4,4,4	0.52	0
65	SPD	3	3234	-	9,9,9	0.34	0	8,8,8	0.78	0
65	SPD	3	3239	-	9,9,9	0.32	0	8,8,8	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
63	PUT	5	1604	-	5,5,5	0.23	0	4,4,4	0.52	0
65	SPD	3	3242	-	9,9,9	0.32	0	8,8,8	0.85	0
65	SPD	5	1601	-	9,9,9	0.42	0	8,8,8	1.41	2 (25%)
66	N2P	5	1602	-	6,6,6	0.24	0	5,5,5	0.63	0
65	SPD	3	3238	-	9,9,9	0.32	0	8,8,8	0.89	0
65	SPD	3	3247	-	9,9,9	0.32	0	8,8,8	0.89	0
65	SPD	3	3250	-	9,9,9	0.32	0	8,8,8	0.85	0
65	SPD	3	3248	-	9,9,9	0.32	0	8,8,8	0.86	0
64	SPM	3	3251	-	13,13,13	0.35	0	12,12,12	0.90	0
66	N2P	3	3249	-	6,6,6	0.24	0	5,5,5	0.64	0
63	PUT	3	3227	-	5,5,5	0.23	0	4,4,4	0.51	0
65	SPD	3	3237	-	9,9,9	0.32	0	8,8,8	0.86	0
65	SPD	3	3240	-	9,9,9	0.32	0	8,8,8	0.85	0
65	SPD	3	3241	-	9,9,9	0.31	0	8,8,8	0.84	0
65	SPD	3	3236	-	9,9,9	0.32	0	8,8,8	0.78	0
65	SPD	3	3252	-	9,9,9	0.18	0	8,8,8	0.18	0
65	SPD	3	3232	-	9,9,9	0.32	0	8,8,8	0.82	0
63	PUT	3	3226	-	5,5,5	0.23	0	4,4,4	0.52	0
65	SPD	3	3235	-	9,9,9	0.32	0	8,8,8	0.87	0
66	N2P	3	3245	-	6,6,6	0.24	0	5,5,5	0.64	0
66	N2P	3	3246	-	6,6,6	0.24	0	5,5,5	0.63	0
60	CLM	3	3001	-	20,20,20	2.07	2 (10%)	23,27,27	0.87	1 (4%)
65	SPD	3	3231	-	9,9,9	0.32	0	8,8,8	0.89	0
64	SPM	b	303	-	13,13,13	0.17	0	12,12,12	0.31	0
63	PUT	3	3233	-	5,5,5	0.24	0	4,4,4	0.51	0

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
63	PUT	3	3225	-	-	0/3/3/3	-
64	SPM	3	3243	-	-	2/11/11/11	-
67	LYS	8	103	9	-	3/6/7/9	-
63	PUT	3	3228	-	-	0/3/3/3	-
64	SPM	3	3229	-	-	4/11/11/11	-
65	SPD	3	3230	-	-	0/7/7/7	-
63	PUT	3	3244	-	-	0/3/3/3	-
65	SPD	5	1603	-	-	1/7/7/7	-
63	PUT	3	3224	-	-	0/3/3/3	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
65	SPD	3	3234	-	-	1/7/7/7	-
65	SPD	3	3239	-	-	1/7/7/7	-
63	PUT	5	1604	-	-	0/3/3/3	-
65	SPD	3	3242	-	-	0/7/7/7	-
65	SPD	5	1601	-	-	1/7/7/7	-
66	N2P	5	1602	-	-	3/4/4/4	-
65	SPD	3	3238	-	-	0/7/7/7	-
65	SPD	3	3247	-	-	0/7/7/7	-
65	SPD	3	3250	-	-	0/7/7/7	-
65	SPD	3	3248	-	-	1/7/7/7	-
64	SPM	3	3251	-	-	2/11/11/11	-
66	N2P	3	3249	-	-	1/4/4/4	-
63	PUT	3	3227	-	-	1/3/3/3	-
65	SPD	3	3237	-	-	0/7/7/7	-
65	SPD	3	3240	-	-	0/7/7/7	-
65	SPD	3	3241	-	-	2/7/7/7	-
65	SPD	3	3236	-	-	2/7/7/7	-
65	SPD	3	3252	-	-	0/7/7/7	-
65	SPD	3	3232	-	-	0/7/7/7	-
63	PUT	3	3226	-	-	0/3/3/3	-
65	SPD	3	3235	-	-	0/7/7/7	-
66	N2P	3	3245	-	-	1/4/4/4	-
66	N2P	3	3246	-	-	2/4/4/4	-
60	CLM	3	3001	-	-	3/20/22/22	0/1/1/1
65	SPD	3	3231	-	-	1/7/7/7	-
64	SPM	b	303	-	-	2/11/11/11	-
63	PUT	3	3233	-	-	0/3/3/3	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
60	3	3001	CLM	C2-N2	7.39	1.49	1.34
60	3	3001	CLM	O9B-N9	-2.64	1.18	1.22

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
65	5	1601	SPD	C4-C3-C2	-2.28	97.82	113.42
60	3	3001	CLM	C3-N2-C2	-2.20	119.42	123.25
65	5	1601	SPD	C4-C5-N6	-2.05	106.56	112.07

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

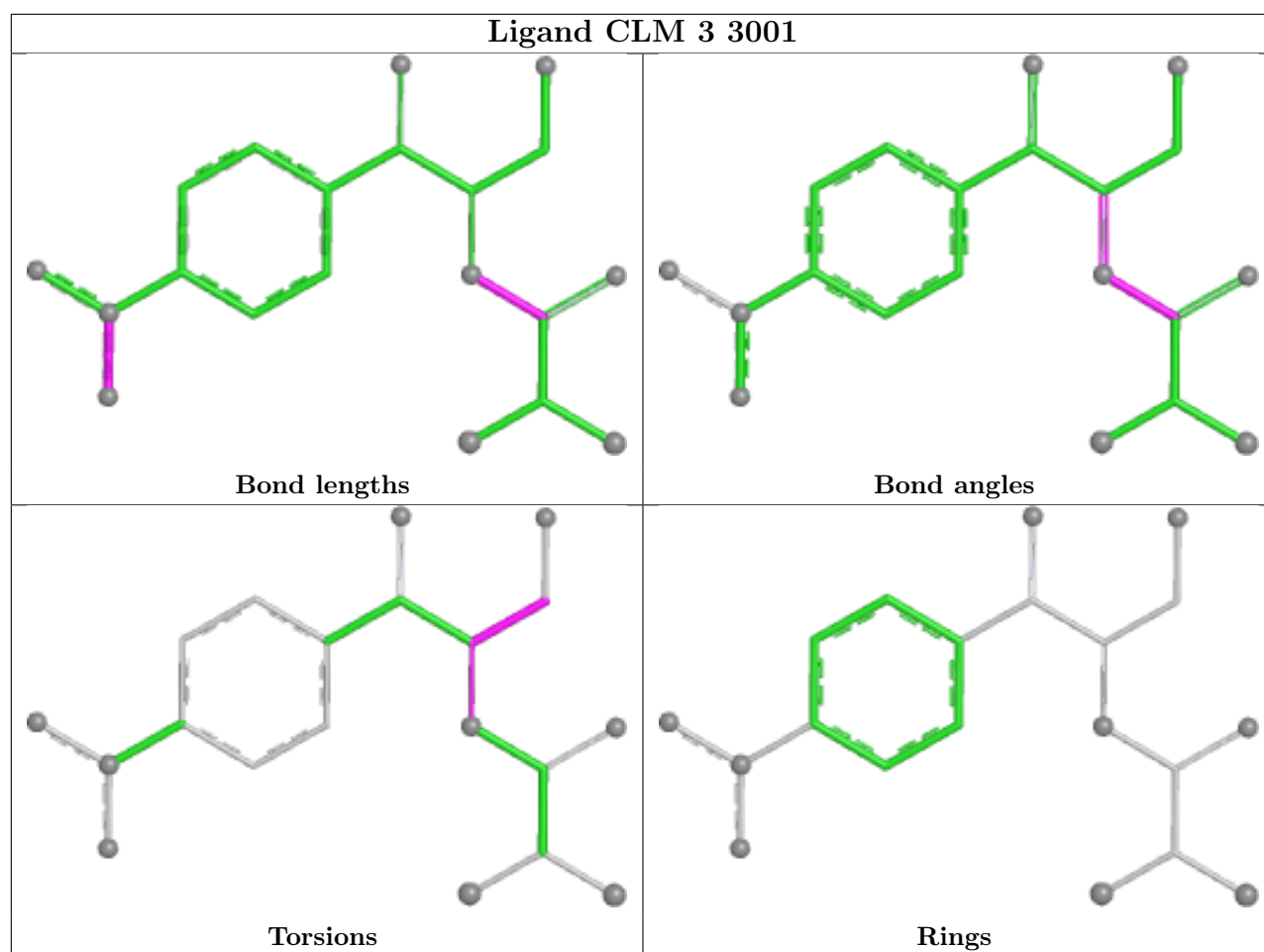
Mol	Chain	Res	Type	Atoms
67	8	103	LYS	O-C-CA-CB
67	8	103	LYS	N-CA-CB-CG
67	8	103	LYS	C-CA-CB-CG
65	5	1603	SPD	C8-C7-N6-C5
66	5	1602	N2P	C2-C3-C4-C5

There are no ring outliers.

12 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
64	3	3243	SPM	2	0
63	3	3228	PUT	1	0
64	3	3229	SPM	4	0
65	5	1603	SPD	1	0
65	5	1601	SPD	6	0
65	3	3250	SPD	1	0
64	3	3251	SPM	1	0
65	3	3241	SPD	1	0
65	3	3252	SPD	1	0
60	3	3001	CLM	1	0
64	b	303	SPM	3	0
63	3	3233	PUT	1	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.

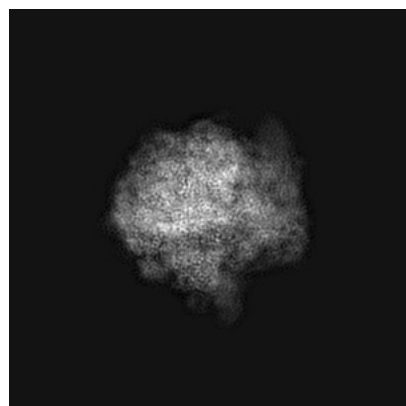
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17135. These allow visual inspection of the internal detail of the map and identification of artifacts.

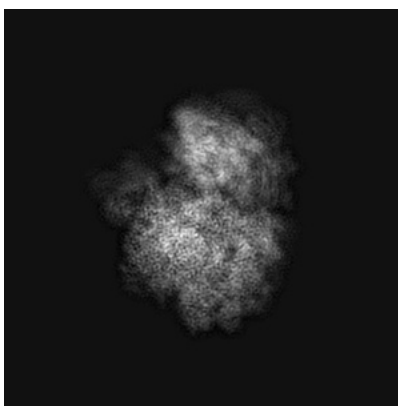
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

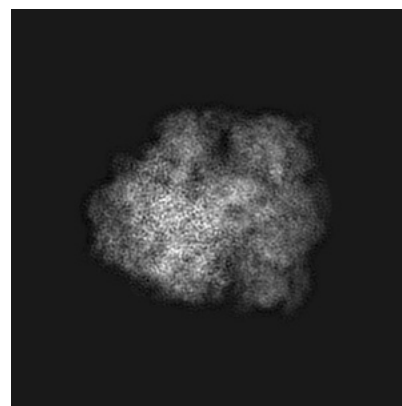
6.1.1 Primary map



X

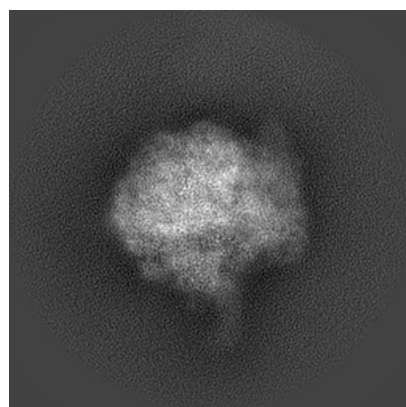


Y

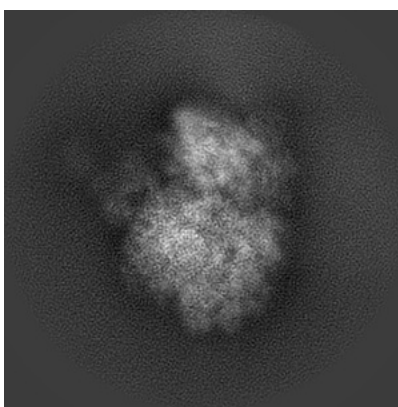


Z

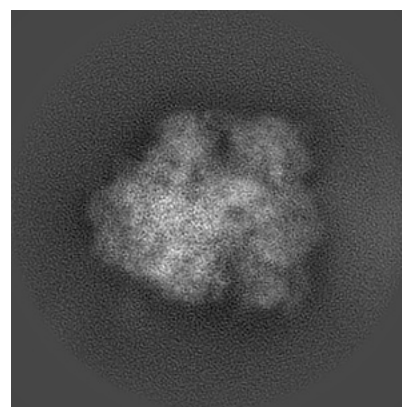
6.1.2 Raw map



X



Y

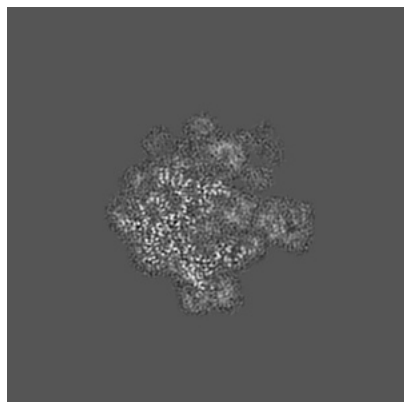


Z

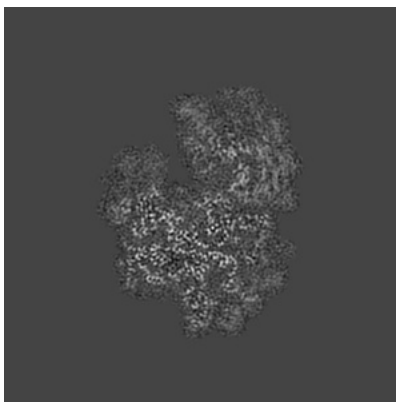
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

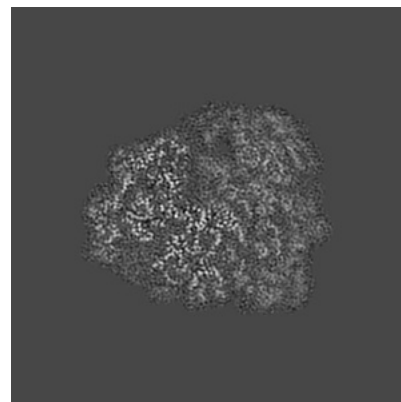
6.2.1 Primary map



X Index: 128

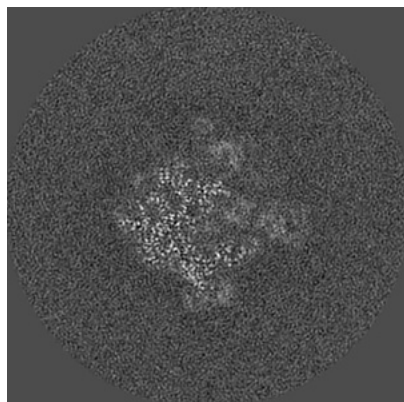


Y Index: 128

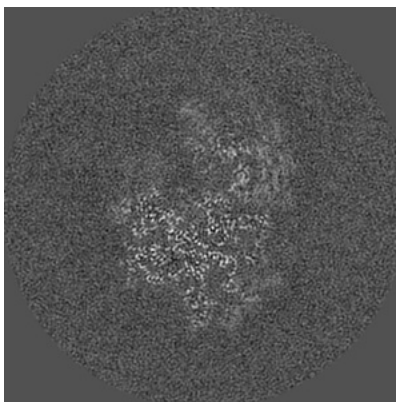


Z Index: 128

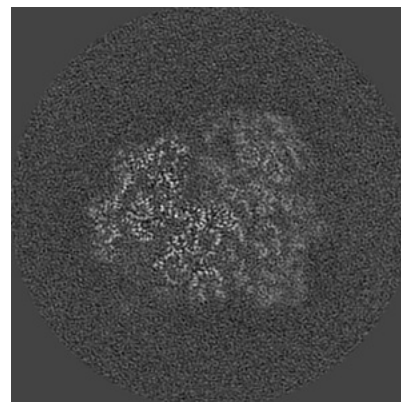
6.2.2 Raw map



X Index: 128



Y Index: 128

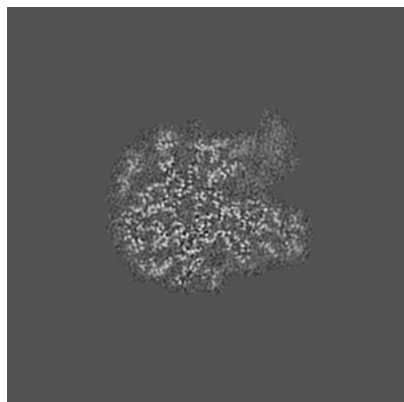


Z Index: 128

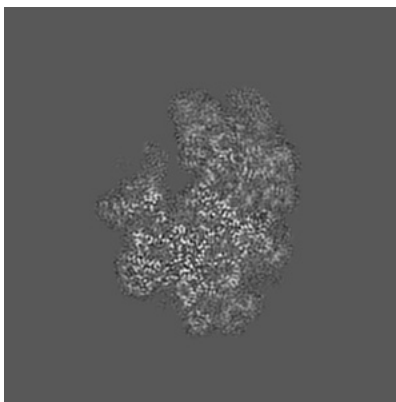
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

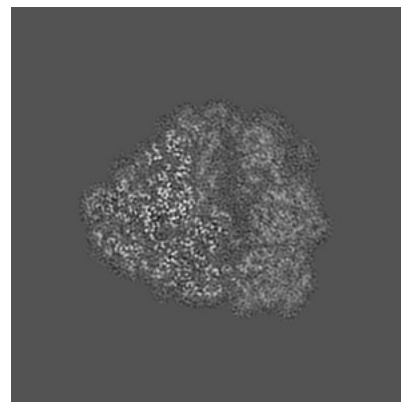
6.3.1 Primary map



X Index: 108

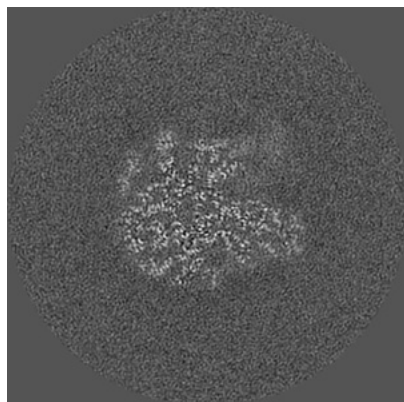


Y Index: 122

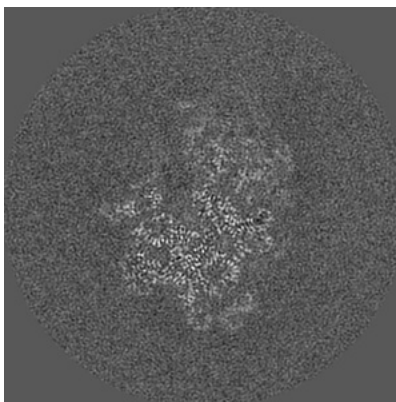


Z Index: 121

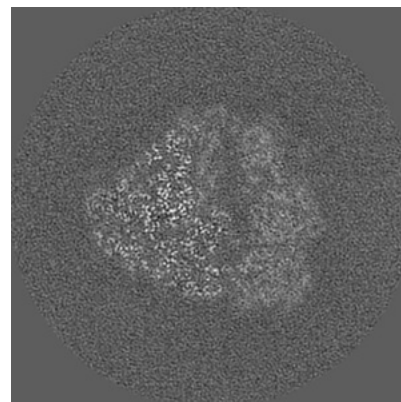
6.3.2 Raw map



X Index: 108



Y Index: 123

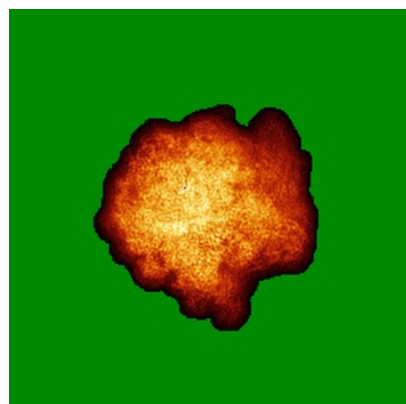


Z Index: 121

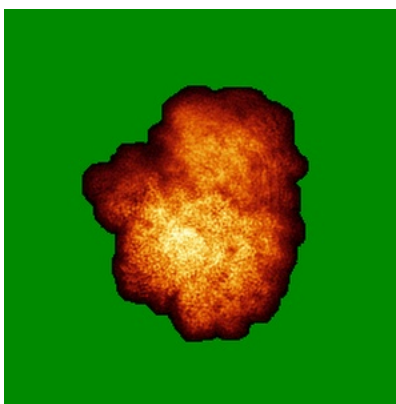
The images above show the largest variance slices of the map in three orthogonal directions.

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

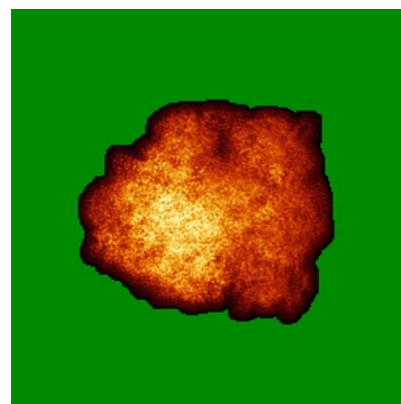
6.4.1 Primary map



X

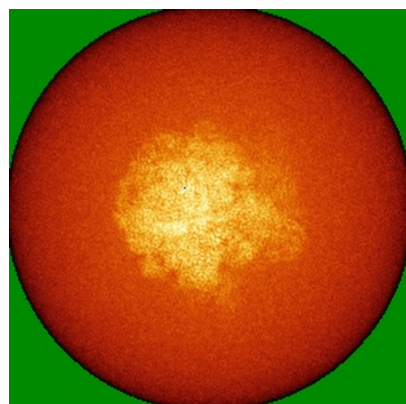


Y

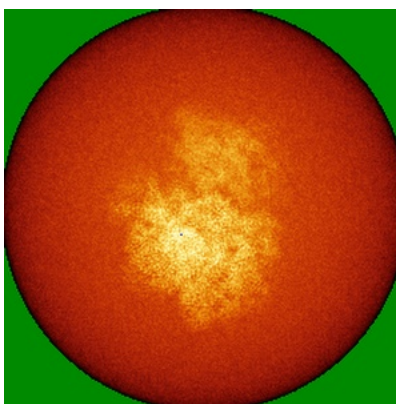


Z

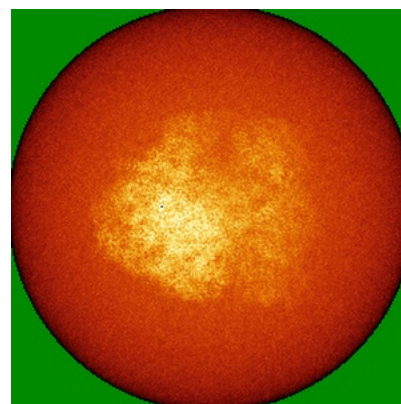
6.4.2 Raw map



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



X



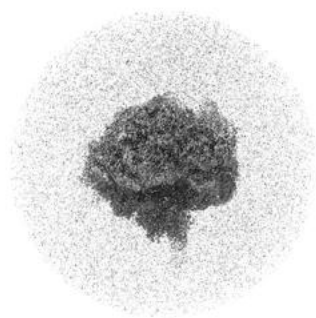
Y



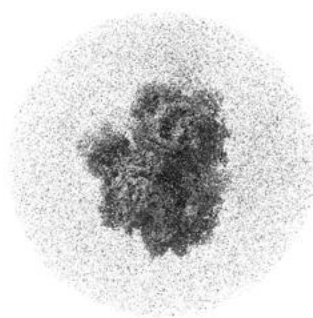
Z

The images above show the 3D surface view of the map at the recommended contour level 0.99. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

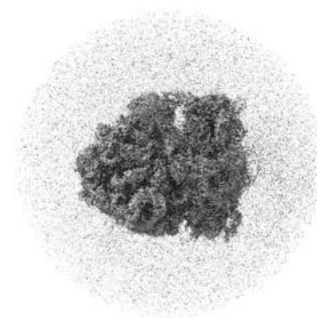
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

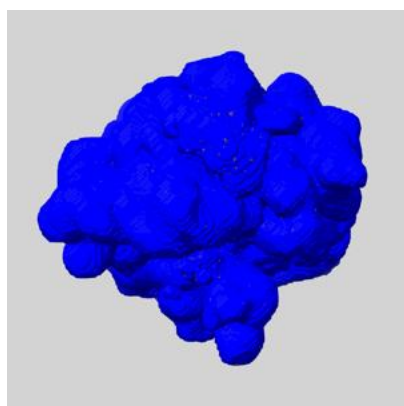
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

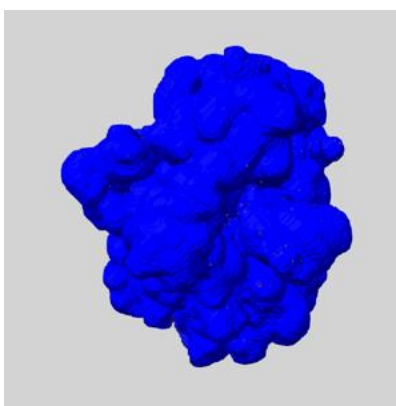
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

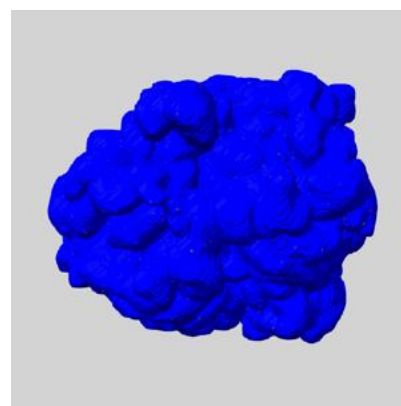
6.6.1 emd_17135_msk_1.map [i](#)



X



Y

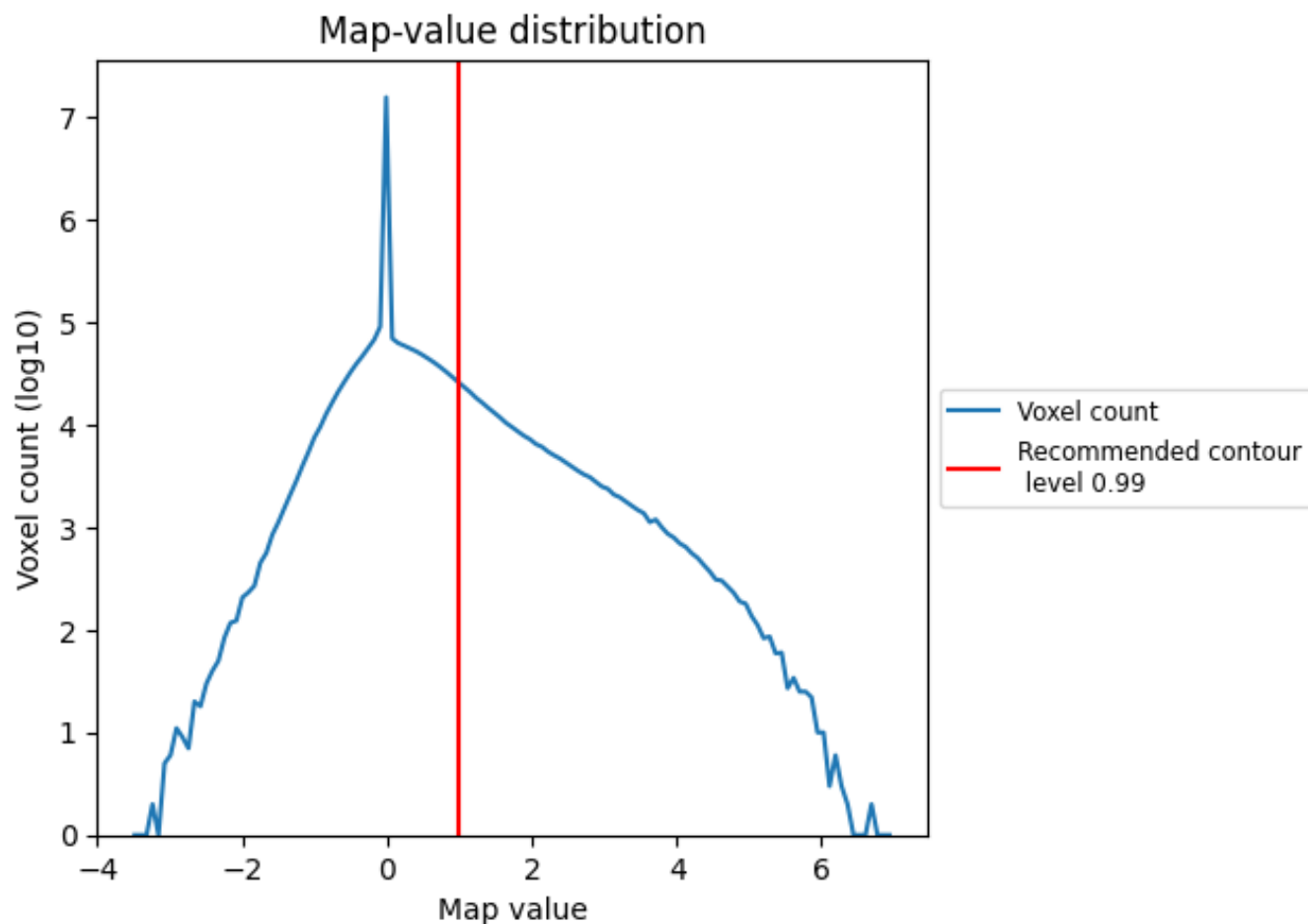


Z

7 Map analysis [i](#)

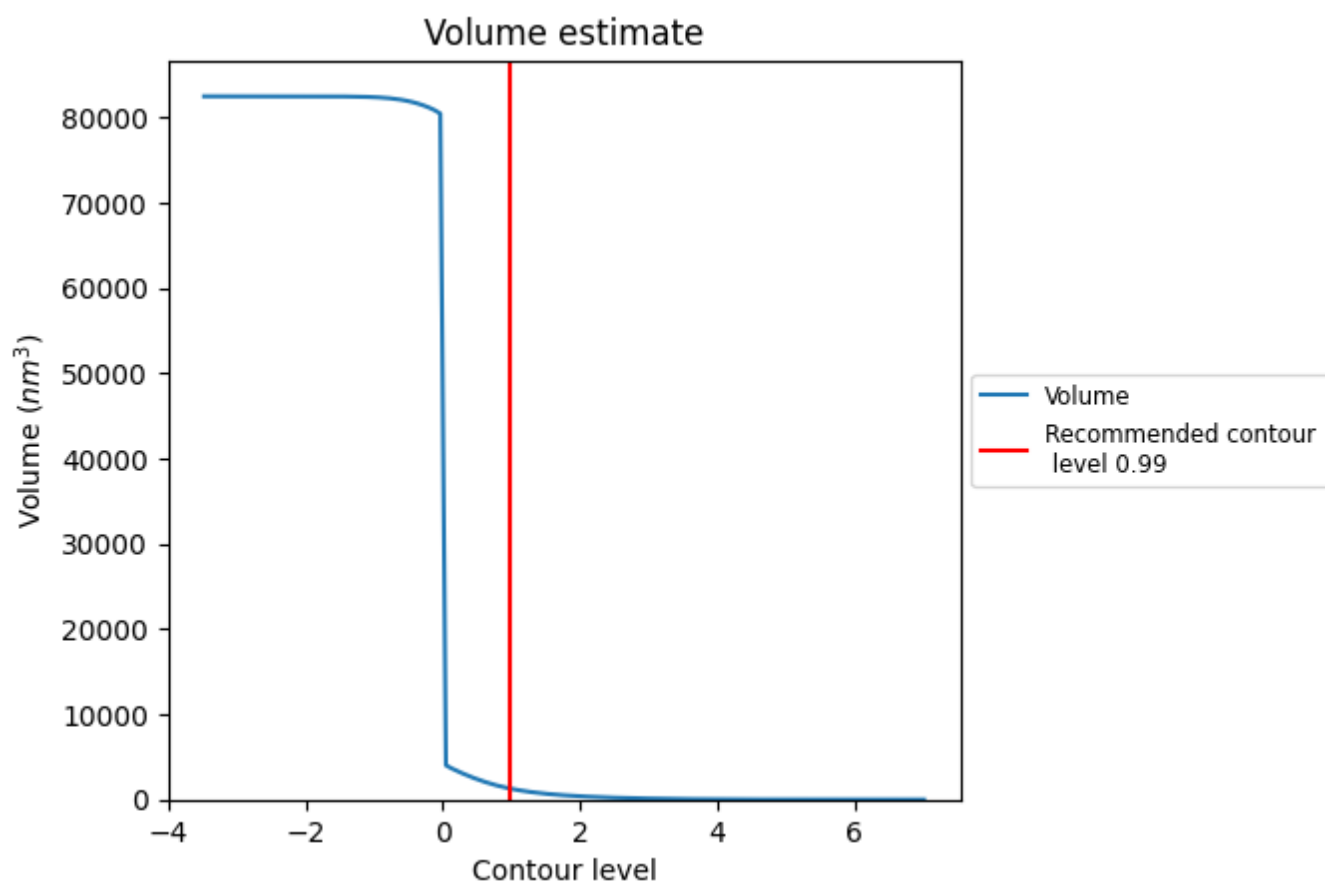
This section contains the results of statistical analysis of the map.

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.

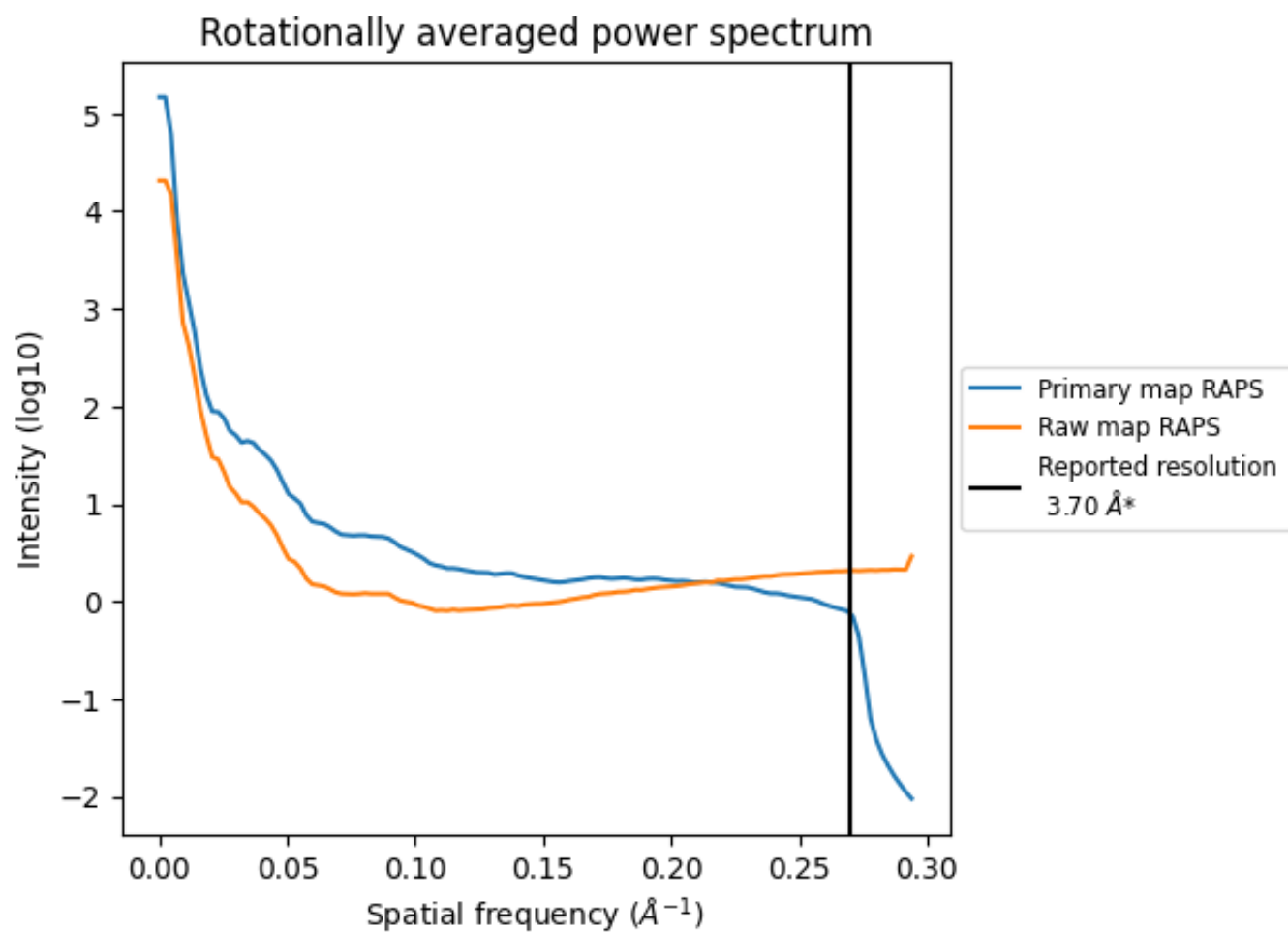
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1276 nm³; this corresponds to an approximate mass of 1153 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 ⓘ

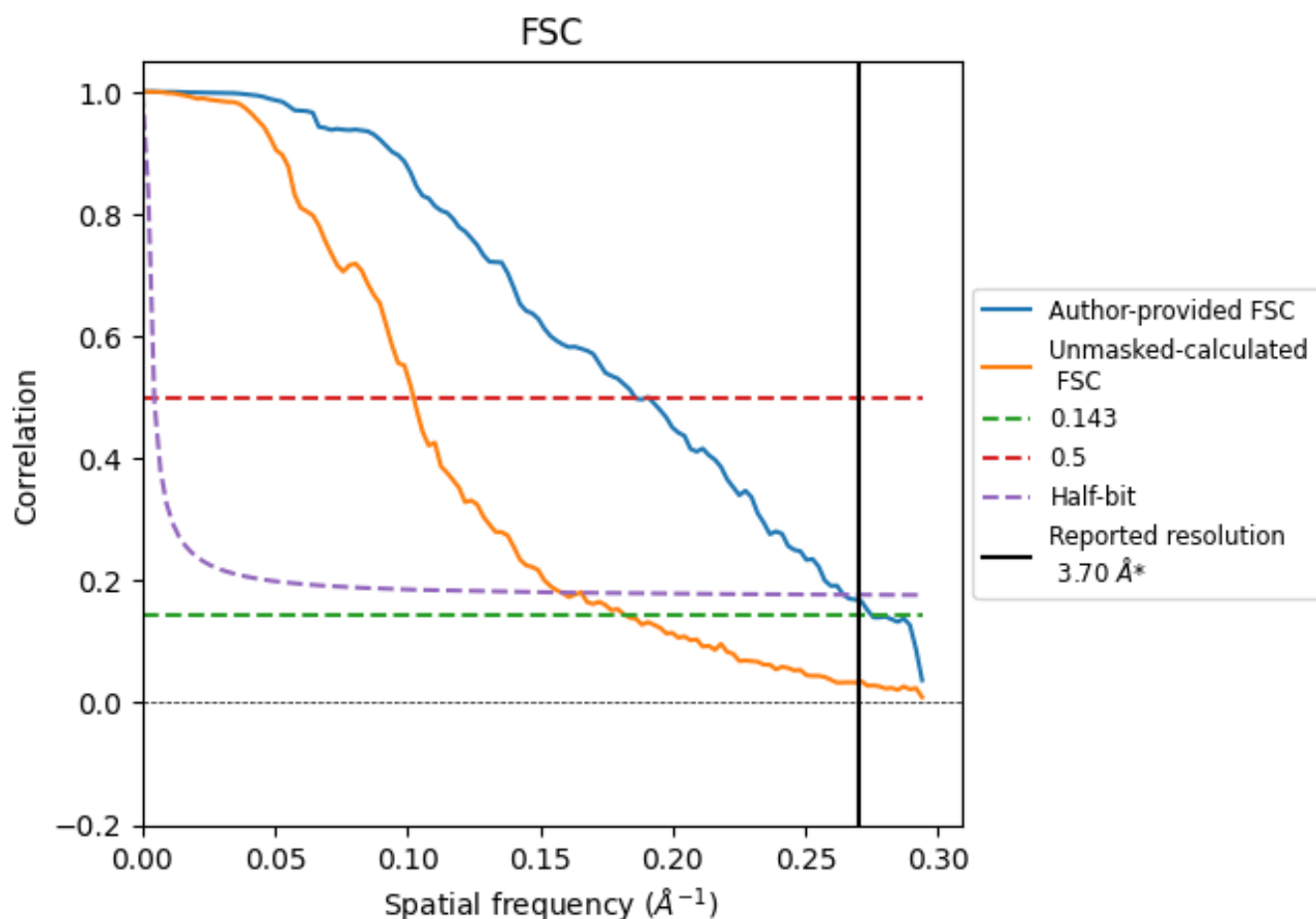


*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

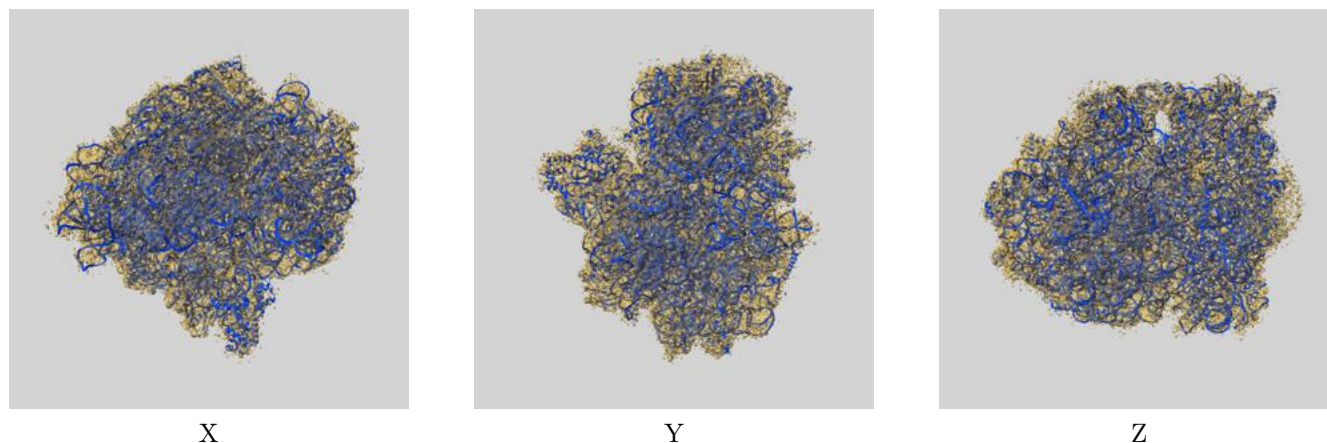
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	3.64	5.39	3.78
Unmasked-calculated*	5.49	9.78	6.32

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 5.49 differs from the reported value 3.7 by more than 10 %

9 Map-model fit [i](#)

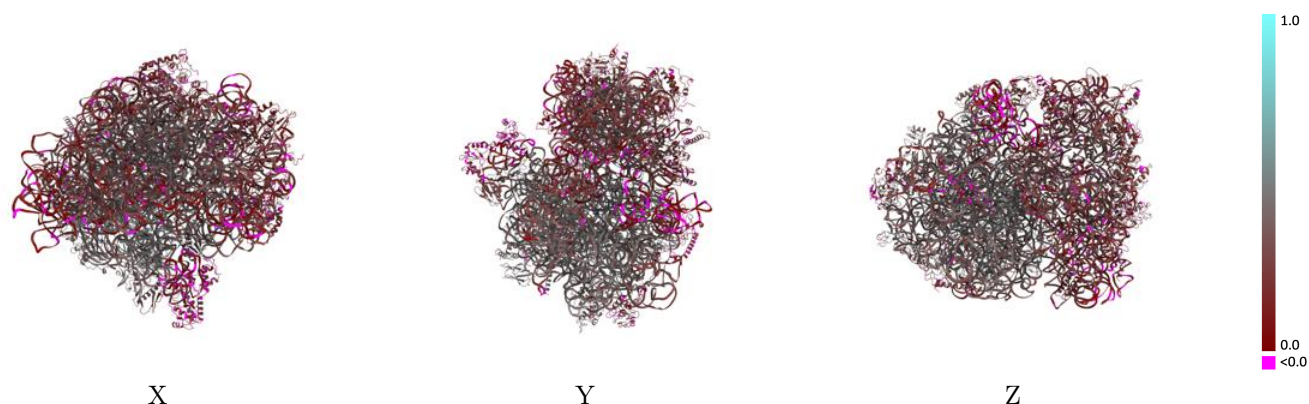
This section contains information regarding the fit between EMDB map EMD-17135 and PDB model 8P7Y. Per-residue inclusion information can be found in section [3](#) on page [21](#).

9.1 Map-model overlay [i](#)



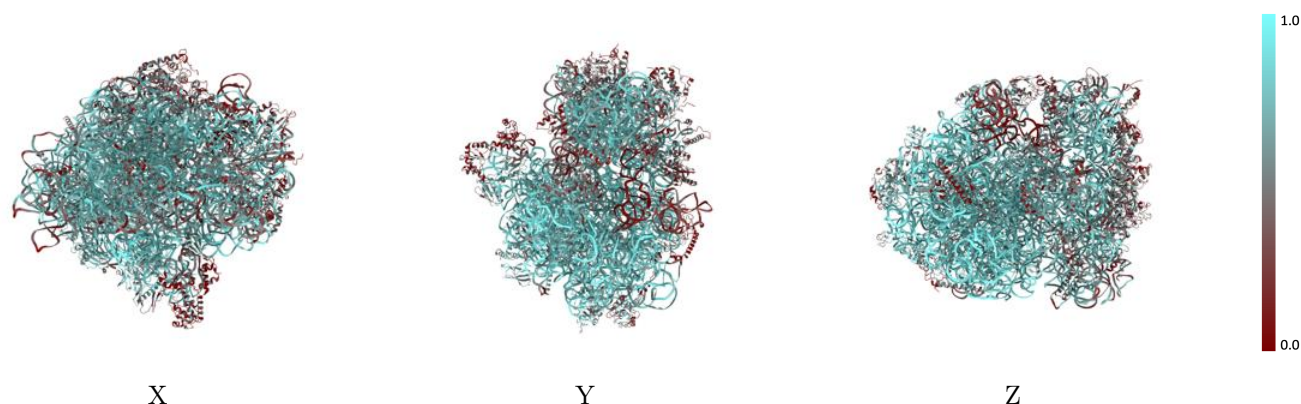
The images above show the 3D surface view of the map at the recommended contour level 0.99 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



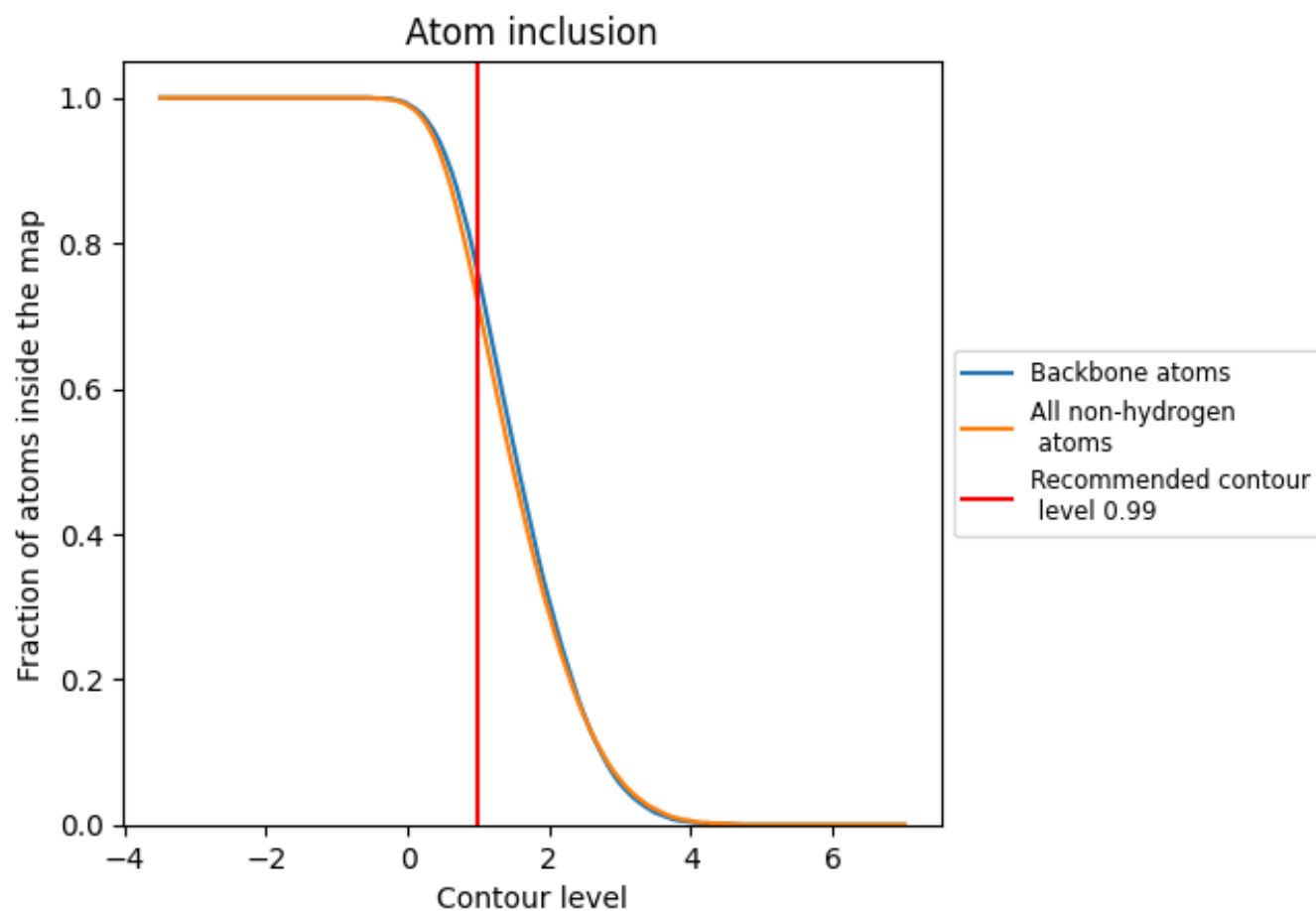
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.99).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 72% 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 (0.99) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7210	 0.3300
0	 0.8210	 0.4660
1	 0.7530	 0.4350
2	 0.7910	 0.4480
3	 0.8630	 0.3880
4	 0.8120	 0.3110
5	 0.7630	 0.2720
6	 0.0860	 0.0460
7	 0.7240	 0.2390
8	 0.4210	 0.2310
A	 0.3900	 0.2460
B	 0.4850	 0.2990
C	 0.3950	 0.2110
D	 0.5440	 0.3410
E	 0.3100	 0.1970
F	 0.3790	 0.2120
G	 0.4650	 0.2610
H	 0.4400	 0.2320
I	 0.4080	 0.2670
J	 0.4710	 0.2480
K	 0.5850	 0.3300
L	 0.3610	 0.1790
M	 0.5700	 0.2930
N	 0.4780	 0.2240
O	 0.4480	 0.1960
P	 0.3730	 0.1870
Q	 0.4190	 0.2090
R	 0.3380	 0.1840
S	 0.5290	 0.2280
T	 0.5040	 0.2900
U	 0.3900	 0.2050
X	 0.1210	 0.1690
Y	 0.5180	 0.2230
Z	 0.2730	 0.1910
a	 0.7740	 0.4350



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Chain	Atom inclusion	Q-score
b	 0.7490	 0.4380
c	 0.7160	 0.4150
d	 0.5080	 0.2690
e	 0.6090	 0.3660
f	 0.1700	 0.1450
g	 0.2200	 0.1100
h	 0.1010	 0.0720
i	 0.7720	 0.4440
j	 0.7270	 0.4500
k	 0.7210	 0.4220
l	 0.7490	 0.4360
m	 0.7840	 0.4510
n	 0.6730	 0.3640
o	 0.6900	 0.4130
p	 0.7640	 0.4330
q	 0.7370	 0.4480
r	 0.7650	 0.4370
s	 0.7080	 0.3990
t	 0.5520	 0.3490
u	 0.7400	 0.4280
v	 0.7610	 0.4340
w	 0.5780	 0.3550
x	 0.2470	 0.1830
y	 0.8120	 0.4620
z	 0.7460	 0.4420