



wwPDB EM Validation Summary Report ⓘ

Jun 21, 2026 – 08:56 am BST

PDB ID : 7ZJW / pdb_00007zjw
EMDB ID : EMD-14751
Title : Rabbit 80S ribosome as it decodes the Sec-UGA codon
Authors : Hilal, T.; Simonovic, M.; Spahn, C.M.T.
Deposited on : 2022-04-12
Resolution : 2.80 Å(reported)
Based on initial model : 7O7Y

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:

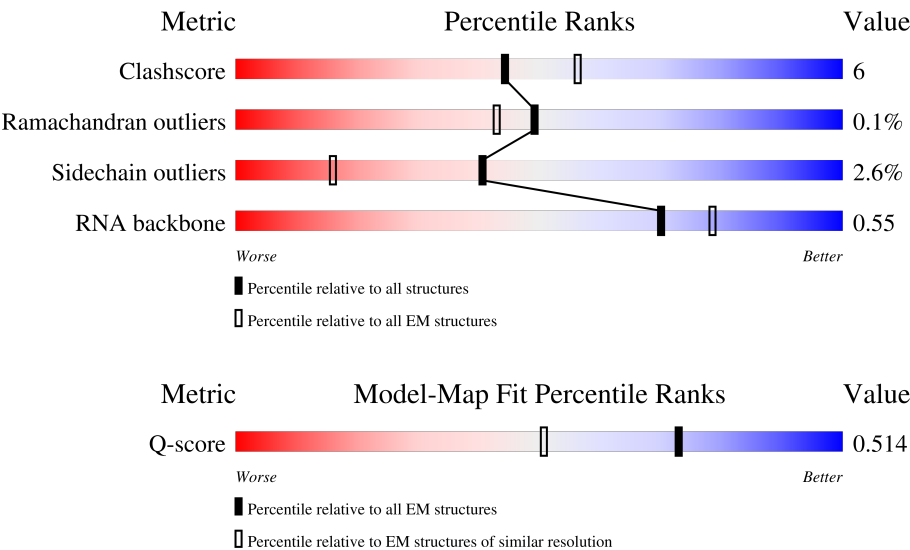
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



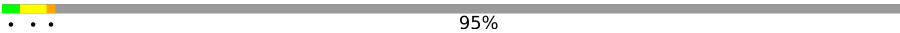
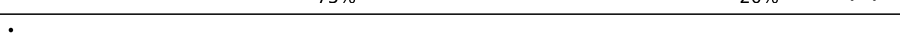
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11806 (2.30 - 3.30)

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 <=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	E	616	8% 71% 13% 16%
2	F	90	10% 44% 49% 7%
3	I	1072	5% 8% 83%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	S	855	 95%
5	B	854	 19% 5% 76%
6	L5	4808	 53% 21% 22%
7	L7	119	 82% 17%
8	L8	158	 78% 20%
9	LD	257	 81% 18%
10	LE	403	 79% 18%
11	LF	413	 75% 12% 12%
12	LG	297	 81% 17%
13	LH	291	 65% 14% 20%
14	LI	247	 74% 17% 9%
15	LJ	266	 77% 10% 12%
16	LK	192	 82% 16%
17	LL	214	 5% 80% 17%
18	LM	178	 75% 20%
19	LO	211	 85% 13%
20	LP	218	 55% 9% 37%
21	LQ	204	 80% 19%
22	LR	203	 87% 11%
23	LS	184	 67% 19% 14%
24	LT	188	 86% 13%
25	LU	196	 73% 15% 8%
26	LV	176	 83% 17%
27	LW	160	 84% 14%
28	LX	128	 62% 15% 23%

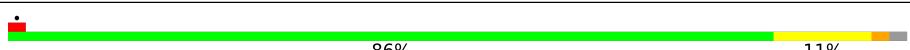
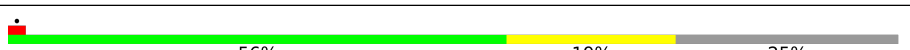
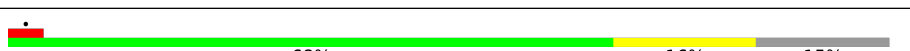
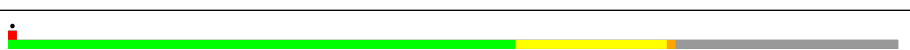
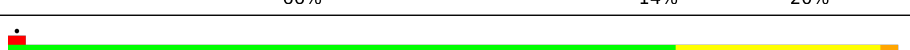
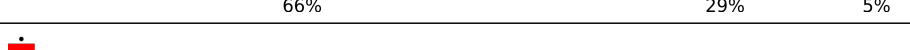
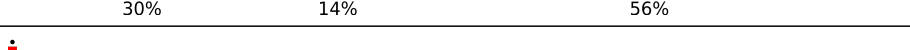
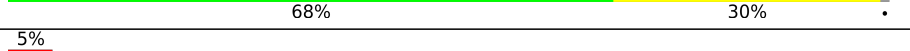
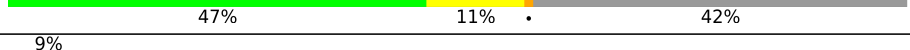
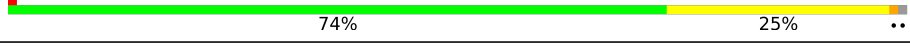
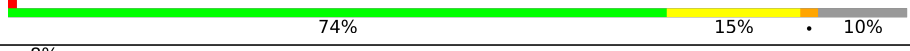
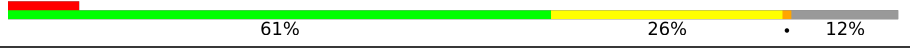
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
29	LY	140	
30	LZ	157	
31	La	156	
32	Lb	145	
33	Lc	136	
34	Ld	148	
35	Le	245	
36	Lf	115	
37	Lg	125	
38	Lh	135	
39	Li	110	
40	Lj	117	
41	Lk	123	
42	Ll	105	
43	Lm	97	
44	Ln	70	
45	Lo	51	
46	Lp	128	
47	Lq	106	
48	Lr	92	
49	Ls	137	
50	Lt	318	
51	Lx	217	
52	S2	1870	
53	SB	84	

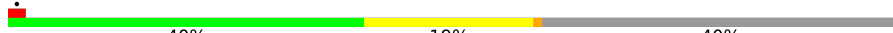
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
54	SC	69	
55	SD	156	
56	SE	133	
57	SF	115	
58	SG	317	
59	SH	56	
60	SL	295	
61	SM	264	
62	SN	293	
63	SO	281	
64	SP	263	
65	SQ	204	
66	SR	249	
67	SS	432	
68	ST	208	
69	SU	194	
70	SV	165	
71	SW	158	
72	SX	132	
73	SY	151	
74	SZ	151	
75	Sa	145	
76	Sb	172	
77	Sc	135	
78	Sd	152	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
79	Se	145	
80	Sf	119	
81	Sg	83	
82	Sh	130	
83	Si	143	
84	Sj	130	
85	Sk	124	
86	Sl	25	

2 Entry composition

There are 93 unique types of molecules in this entry. The entry contains 230266 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 eEFSec.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	E	519	Total	C	N	O	S	0	0
			3335	2110	616	599	10		

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-19	MET	-	initiating methionine	UNP P57772
E	-18	GLY	-	expression tag	UNP P57772
E	-17	SER	-	expression tag	UNP P57772
E	-16	SER	-	expression tag	UNP P57772
E	-15	HIS	-	expression tag	UNP P57772
E	-14	HIS	-	expression tag	UNP P57772
E	-13	HIS	-	expression tag	UNP P57772
E	-12	HIS	-	expression tag	UNP P57772
E	-11	HIS	-	expression tag	UNP P57772
E	-10	HIS	-	expression tag	UNP P57772
E	-9	SER	-	expression tag	UNP P57772
E	-8	SER	-	expression tag	UNP P57772
E	-7	GLY	-	expression tag	UNP P57772
E	-6	LEU	-	expression tag	UNP P57772
E	-5	VAL	-	expression tag	UNP P57772
E	-4	PRO	-	expression tag	UNP P57772
E	-3	ARG	-	expression tag	UNP P57772
E	-2	GLY	-	expression tag	UNP P57772
E	-1	SER	-	expression tag	UNP P57772
E	0	HIS	-	expression tag	UNP P57772
E	409	GLU	GLN	conflict	UNP P57772

- Molecule 2 is a RNA chain called human Ser-tRNA-Sec.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	F	90	Total	C	N	O	P	0	0
			1914	853	334	637	90		

- Molecule 3 is a RNA chain called CrPV IRES.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	183	Total	C	N	O	P	0	0
			3882	1739	675	1285	183		

- Molecule 4 is a RNA chain called GPX4 SECIS element.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	S	46	Total	C	N	O	P	0	0
			983	438	179	320	46		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	1118	U	C	conflict	GB 25123295
S	1126	A	C	conflict	GB 25123295
S	1132	A	U	conflict	GB 25123295
S	1133	G	-	insertion	GB 25123295
S	1134	C	-	insertion	GB 25123295
S	1135	C	-	insertion	GB 25123295
S	1136	C	-	insertion	GB 25123295
S	1150	U	A	conflict	GB 25123295

- Molecule 5 is a protein called SECIS Binding Protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	B	204	Total	C	N	O	S	0	0
			1379	881	257	238	3		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	687	ARG	LYS	conflict	UNP Q96T21
B	692	ILE	VAL	conflict	UNP Q96T21

- Molecule 6 is a RNA chain called 28S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L5	3764	Total	C	N	O	P	0	0
			80771	36002	14762	26243	3764		

- Molecule 7 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	L7	119	Total	C	N	O	P	0	0
			2538	1131	451	837	119		

- Molecule 8 is a RNA chain called 5.8S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	L8	156	Total	C	N	O	P	0	0
			3319	1481	585	1097	156		

- Molecule 9 is a protein called 60S ribosomal protein uL2.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	LD	253	Total	C	N	O	S	0	0
			1939	1214	396	323	6		

- Molecule 10 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	LE	398	Total	C	N	O	S	0	0
			3206	2042	605	546	13		

- Molecule 11 is a protein called 60S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	LF	362	Total	C	N	O	S	0	0
			2886	1814	577	481	14		

- Molecule 12 is a protein called Ribosomal_L18_c domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	LG	294	Total	C	N	O	S	0	0
			2398	1516	439	429	14		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LG	2	AAC	GLY	conflict	UNP G1SYJ6

- Molecule 13 is a protein called 60S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	LH	232	Total	C	N	O	S	0	0
			1859	1197	354	305	3		

- Molecule 14 is a protein called 60S ribosomal Protein uL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	LI	226	Total	C	N	O	S	0	0
			1886	1211	362	304	9		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LI	182	ASN	GLY	conflict	UNP A0A7J8C453
LI	199	HIS	ARG	conflict	UNP A0A7J8C453

- Molecule 15 is a protein called 60S ribosomal protein eL8.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	LJ	233	Total	C	N	O	S	0	0
			1877	1197	361	315	4		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LJ	184	LEU	ILE	conflict	UNP P62424

- Molecule 16 is a protein called 60S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	LK	190	Total	C	N	O	S	0	0
			1516	954	284	272	6		

- Molecule 17 is a protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	LL	213	Total	C	N	O	S	0	0
			1717	1086	332	285	14		

- Molecule 18 is a protein called 60S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	LM	170	Total	C	N	O	S	0	0
			1362	861	254	241	6		

- Molecule 19 is a protein called 60S ribosomal protein eL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	LO	210	Total	C	N	O	S	0	0
			1702	1065	354	279	4		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LO	74	ARG	HIS	conflict	UNP G1TKB3
LO	190	ARG	HIS	conflict	UNP G1TKB3

- Molecule 20 is a protein called 60S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	LP	138	Total	C	N	O	S	0	0
			1137	727	221	182	7		

- Molecule 21 is a protein called 60S ribosomal protein L15.

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

- Molecule 22 is a protein called 60S ribosomal protein uL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	LR	199	Total	C	N	O	S	0	0
			1630	1051	319	255	5		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LR	174	LEU	ILE	conflict	UNP A0A0N8ETI8
LR	194	ASP	GLU	conflict	UNP A0A0N8ETI8

- Molecule 23 is a protein called 60S ribosomal protein uL22.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	LS	159	Total	C	N	O	S	0	0
			1288	808	249	222	9		

- Molecule 24 is a protein called 60S ribosomal Protein eL18.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	LT	187	Total	C	N	O	S	0	0
			1515	946	315	250	4		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LT	134	ARG	CYS	conflict	UNP F6QKI9

- Molecule 25 is a protein called 60S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	LU	180	Total	C	N	O	S	0	0
			1508	933	328	238	9		

- Molecule 26 is a protein called 60S ribosomal protein eL20.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	LV	176	Total	C	N	O	S	0	0
			1457	924	288	234	11		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LV	36	ASN	ILE	conflict	UNP A0A1Z5LHJ5

- Molecule 27 is a protein called 60S ribosomal protein eL21.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	LW	159	Total	C	N	O	S	0	0
			1298	823	252	217	6		

- Molecule 28 is a protein called 60S ribosomal protein eL22.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	LX	99	Total	C	N	O	S	0	0
			806	516	141	147	2		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LX	60	ALA	VAL	conflict	UNP Q4R5I3

- Molecule 29 is a protein called 60S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	LY	139	Total	C	N	O	S	0	0
			1034	648	199	182	5		

- Molecule 30 is a protein called 60S ribosomal protein eL24.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	LZ	93	Total	C	N	O	S	0	0
			766	480	153	129	4		

- Molecule 31 is a protein called 60S ribosomal protein uL23.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	La	118	Total	C	N	O	S	0	0
			967	618	181	167	1		

- Molecule 32 is a protein called 60S ribosomal protein L26.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	Lb	134	Total	C	N	O	S	0	0
			1115	700	226	186	3		

- Molecule 33 is a protein called 60S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	Lc	135	Total	C	N	O	S	0	0
			1107	714	208	182	3		

- Molecule 34 is a protein called 60S ribosomal protein L27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	Ld	147	Total	C	N	O	S	0	0
			1162	734	239	185	4		

- Molecule 35 is a protein called 60S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	Le	108	Total	C	N	O	S	0	0
			881	548	196	134	3		

- Molecule 36 is a protein called 60S ribosomal protein eL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	Lf	108	Total	C	N	O	S	0	0
			836	530	148	151	7		

- Molecule 37 is a protein called 60S ribosomal protein eL31.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	Lg	107	Total	C	N	O	S	0	0
			888	560	171	155	2		

- Molecule 38 is a protein called 60S ribosomal protein eL32.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	Lh	130	Total	C	N	O	S	0	0
			1070	676	221	168	5		

- Molecule 39 is a protein called 60S ribosomal protein eL33.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	Li	110	Total	C	N	O	S	0	0
			884	560	175	144	5		

- Molecule 40 is a protein called 60S ribosomal protein L34.

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

- Molecule 41 is a protein called 60S ribosomal protein uL29.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	Lk	122	Total	C	N	O	S	0	0
			1013	640	204	168	1		

- Molecule 42 is a protein called 60S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	Ll	102	Total	C	N	O	S	0	0
			830	520	176	129	5		

- Molecule 43 is a protein called 60S ribosomal protein L37.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	Lm	86	Total	C	N	O	S	0	0
			705	434	155	111	5		

- Molecule 44 is a protein called 60S ribosomal protein eL38.

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

- Molecule 45 is a protein called 60S ribosomal protein eL39.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	Lo	50	Total	C	N	O	S	0	0
			447	286	96	64	1		

- Molecule 46 is a protein called 60S ribosomal protein L40.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	Lp	52	Total	C	N	O	S	0	0
			431	269	90	66	6		

- Molecule 47 is a protein called 60S ribosomal protein eL42.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	Lq	105	Total	C	N	O	S	0	0
			863	543	175	139	6		

- Molecule 48 is a protein called 60S ribosomal protein eL43.

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

- Molecule 49 is a protein called 60S ribosomal protein eL28.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	Ls	126	Total	C	N	O	S	0	0
			1014	629	209	170	6		

- Molecule 50 is a protein called 60S ribosomal protein P0.

Mol	Chain	Residues	Atoms				AltConf	Trace
50	Lt	115	Total	C	N	O	0	0
			567	337	115	115		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Lt	106	LEU	LYS	conflict	UNP G1SPK4

- Molecule 51 is a protein called 60S ribosomal protein uL1.

Mol	Chain	Residues	Atoms				AltConf	Trace
51	Lx	196	Total	C	N	O	0	0
			972	580	196	196		

- Molecule 52 is a RNA chain called 18S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	S2	1770	Total	C	N	O	P	0	0
			37825	16906	6780	12369	1770		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S2	251	C	U	conflict	GB 37956930
S2	583	U	C	conflict	GB 37956930
S2	584	A	C	conflict	GB 37956930
S2	585	A	G	conflict	GB 37956930
S2	1338	4AC	C	conflict	GB 37956930
S2	1843	4AC	C	conflict	GB 37956930

- Molecule 53 is a protein called 40S ribosomal protein S27.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	SB	83	Total	C	N	O	S	0	0
			651	408	121	115	7		

- Molecule 54 is a protein called 40S ribosomal protein S28.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	SC	63	Total	C	N	O	S	0	0
			495	302	98	93	2		

- Molecule 55 is a protein called 40S ribosomal protein S27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	SD	74	Total	C	N	O	S	0	0
			610	385	117	101	7		

- Molecule 56 is a protein called 40S ribosomal protein S30.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	SE	57	Total	C	N	O	S	0	0
			457	282	101	73	1		

- Molecule 57 is a protein called Ribosomal protein eS26.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	SF	101	Total	C	N	O	S	0	0
			814	507	170	132	5		

- Molecule 58 is a protein called RACK1.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	SG	313	Total	C	N	O	S	0	0
			2436	1535	424	465	12		

- Molecule 59 is a protein called 40S ribosomal protein uS14.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	SH	55	Total	C	N	O	S	0	0
			459	286	94	74	5		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
60	SL	221	Total	C	N	O	S	0	0
			1743	1107	305	323	8		

- Molecule 61 is a protein called 40S ribosomal protein S3a.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	SM	224	Total	C	N	O	S	0	0
			1815	1152	328	321	14		

- Molecule 62 is a protein called 40S ribosomal protein uS5.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	SN	220	Total	C	N	O	S	0	0
			1706	1105	292	300	9		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
SN	33	ILE	VAL	conflict	UNP O18789
SN	101	ALA	SER	conflict	UNP O18789

- Molecule 63 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
63	SO	225	Total	C	N	O	S	0	0
			1751	1116	315	313	7		

- Molecule 64 is a protein called 40S ribosomal protein eS4.

Mol	Chain	Residues	Atoms					AltConf	Trace
64	SP	262	Total	C	N	O	S	0	0
			2076	1324	386	358	8		

- Molecule 65 is a protein called Ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
65	SQ	191	Total	C	N	O	S	0	0
			1509	943	286	273	7		

- Molecule 66 is a protein called 40S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
66	SR	237	Total	C	N	O	S	0	0
			1923	1200	387	329	7		

- Molecule 67 is a protein called 40S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
67	SS	190	Total	C	N	O	S	0	0
			1529	975	281	272	1		

- Molecule 68 is a protein called 40S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
68	ST	206	Total	C	N	O	S	0	0
			1686	1058	332	291	5		

- Molecule 69 is a protein called 40S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
69	SU	185	Total	C	N	O	S	0	0
			1525	969	306	248	2		

- Molecule 70 is a protein called 40S ribosomal protein eS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
70	SV	96	Total	C	N	O	S	0	0
			810	530	143	131	6		

- Molecule 71 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
71	SW	154	Total	C	N	O	S	0	0
			1262	804	236	216	6		

- Molecule 72 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
72	SX	124	Total	C	N	O	S	0	0
			958	600	170	179	9		

- Molecule 73 is a protein called 40S ribosomal protein uS15.

Mol	Chain	Residues	Atoms					AltConf	Trace
73	SY	150	Total	C	N	O	S	0	0
			1208	773	229	205	1		

- Molecule 74 is a protein called 40S ribosomal protein uS11.

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

- Molecule 75 is a protein called 40S ribosomal protein uS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
75	Sa	128	Total	C	N	O	S	0	0
			1048	665	197	179	7		

- Molecule 76 is a protein called 40S ribosomal protein uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
76	Sb	141	Total	C	N	O	S	0	0
			1124	715	212	194	3		

- Molecule 77 is a protein called 40S ribosomal protein eS17.

Mol	Chain	Residues	Atoms					AltConf	Trace
77	Sc	134	Total	C	N	O	S	0	0
			1080	678	201	197	4		

- Molecule 78 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
78	Sd	148	Total	C	N	O	S	0	0
			1217	763	245	208	1		

- Molecule 79 is a protein called 40S Ribosomal protein eS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
79	Se	143	Total	C	N	O	S	0	0
			1113	698	214	198	3		

- Molecule 80 is a protein called 40S ribosomal protein uS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
80	Sf	104	Total	C	N	O	S	0	0
			821	514	155	148	4		

- Molecule 81 is a protein called 40S ribosomal protein eS21.

Mol	Chain	Residues	Atoms					AltConf	Trace
81	Sg	83	Total	C	N	O	S	0	0
			640	394	117	124	5		

- Molecule 82 is a protein called 40S ribosomal protein S15a.

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

- Molecule 83 is a protein called 40S ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
83	Si	141	Total	C	N	O	S	0	0
			1099	693	219	184	3		

- Molecule 84 is a protein called 40S ribosomal protein S24.

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

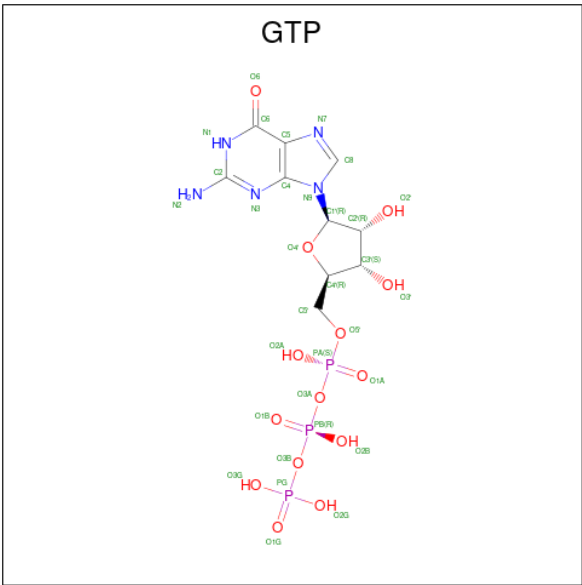
- Molecule 85 is a protein called 40S ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
85	Sk	75	Total	C	N	O	S	0	0
			598	382	111	104	1		

- Molecule 86 is a protein called 60S ribosomal protein L41.

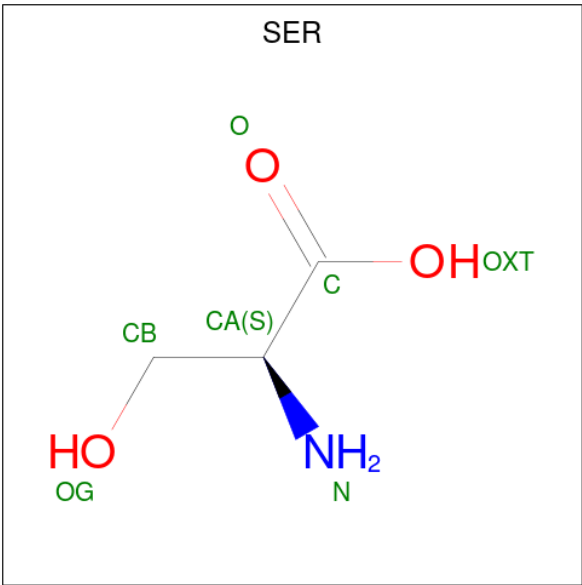
Mol	Chain	Residues	Atoms					AltConf	Trace
86	Sl	25	Total	C	N	O	S	0	0
			239	145	64	27	3		

- Molecule 87 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					AltConf
87	E	1	Total	C	N	O	P	0
			32	10	5	14	3	
87	L7	1	Total	C	N	O	P	0
			32	10	5	14	3	

- Molecule 88 is SERINE (CCD ID: SER) (formula: C₃H₇NO₃).



Mol	Chain	Residues	Atoms				AltConf
88	F	1	Total	C	N	O	0
			6	3	1	2	

- Molecule 89 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
89	I	2	Total 2	Mg 2	0
89	L5	162	Total 162	Mg 162	0
89	L7	2	Total 2	Mg 2	0
89	L8	1	Total 1	Mg 1	0
89	LD	1	Total 1	Mg 1	0
89	LS	1	Total 1	Mg 1	0
89	LY	1	Total 1	Mg 1	0
89	S2	55	Total 55	Mg 55	0

- Molecule 90 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
90	L5	17	Total 17	K 17	0
90	S2	1	Total 1	K 1	0

- Molecule 91 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
91	Lj	1	Total 1	Zn 1	0
91	Lm	1	Total 1	Zn 1	0
91	Lp	1	Total 1	Zn 1	0
91	Lq	1	Total 1	Zn 1	0
91	Lr	1	Total 1	Zn 1	0
91	SD	1	Total 1	Zn 1	0
91	SF	2	Total 2	Zn 2	0
91	SH	1	Total 1	Zn 1	0

- Molecule 92 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		AltConf
92	SZ	1	Total 1	Na 1	0

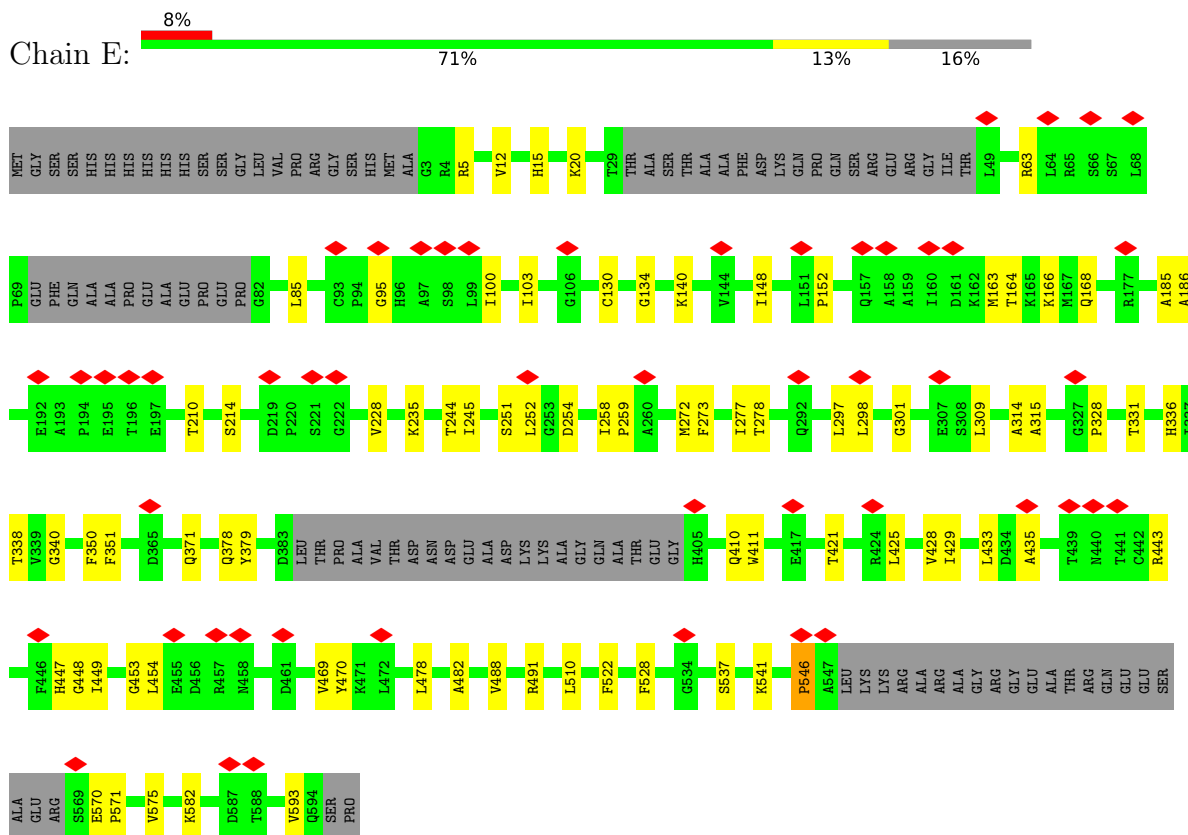
- Molecule 93 is water.

Mol	Chain	Residues	Atoms		AltConf
93	L5	56	Total 56	O 56	0
93	L7	1	Total 1	O 1	0
93	L8	1	Total 1	O 1	0
93	LD	1	Total 1	O 1	0
93	LE	1	Total 1	O 1	0
93	LH	1	Total 1	O 1	0
93	LK	1	Total 1	O 1	0
93	LL	1	Total 1	O 1	0
93	LQ	1	Total 1	O 1	0
93	LV	1	Total 1	O 1	0
93	Ld	1	Total 1	O 1	0
93	S2	34	Total 34	O 34	0
93	SF	1	Total 1	O 1	0
93	SH	1	Total 1	O 1	0
93	SW	1	Total 1	O 1	0
93	Sd	1	Total 1	O 1	0
93	Se	1	Total 1	O 1	0
93	Si	2	Total 2	O 2	0

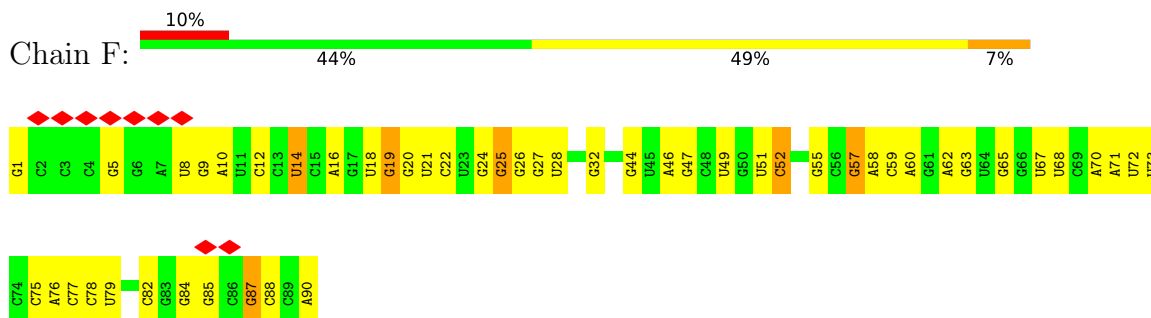
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: eEFSec



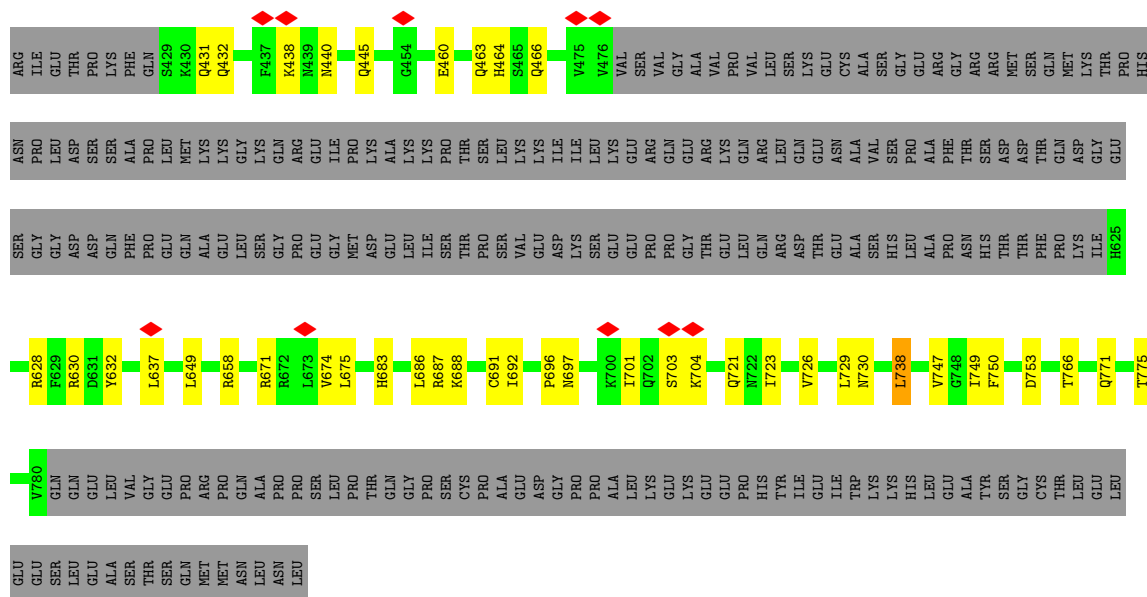
- Molecule 2: human Ser-tRNA-Sec



- Molecule 3: CrPV IRES

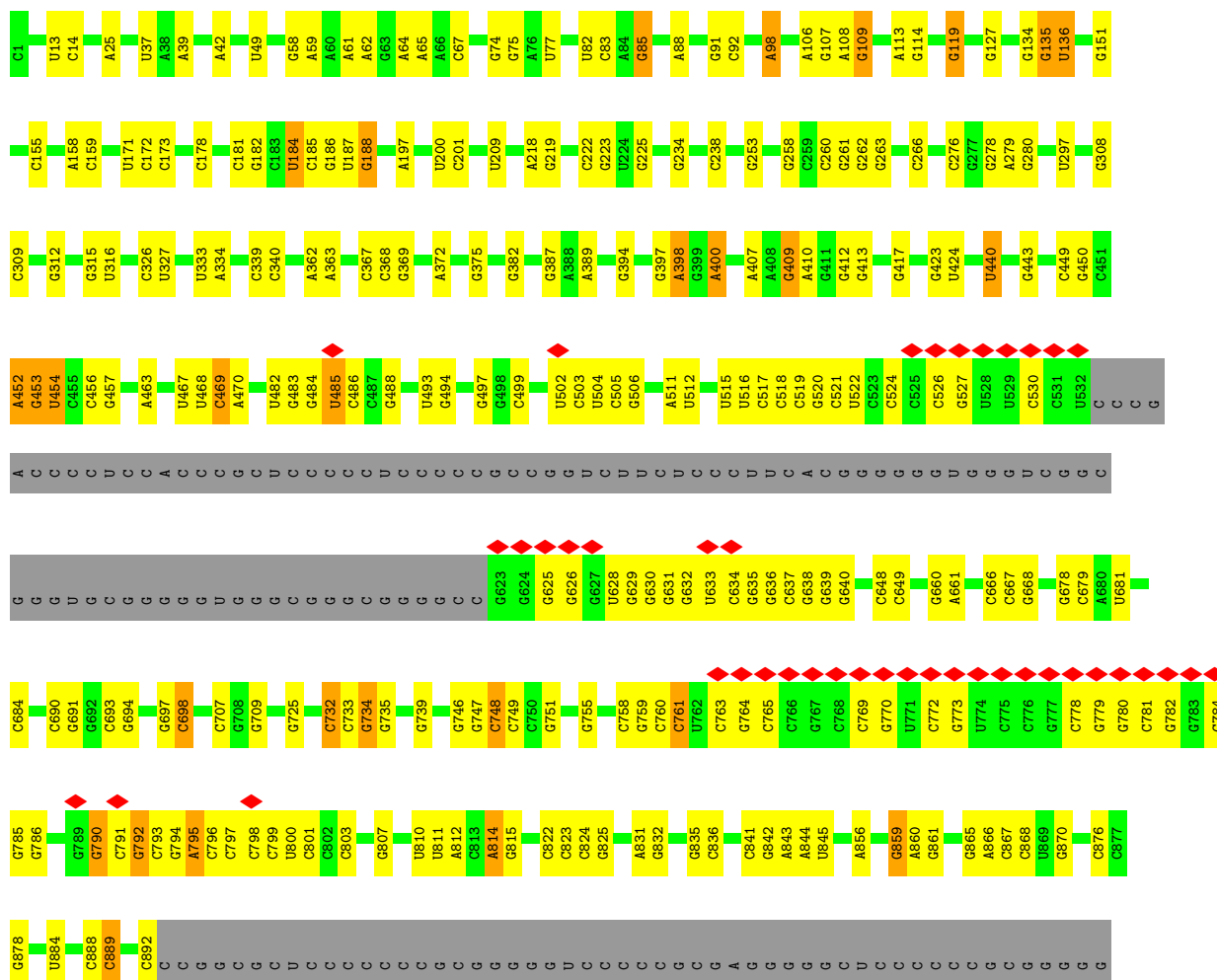


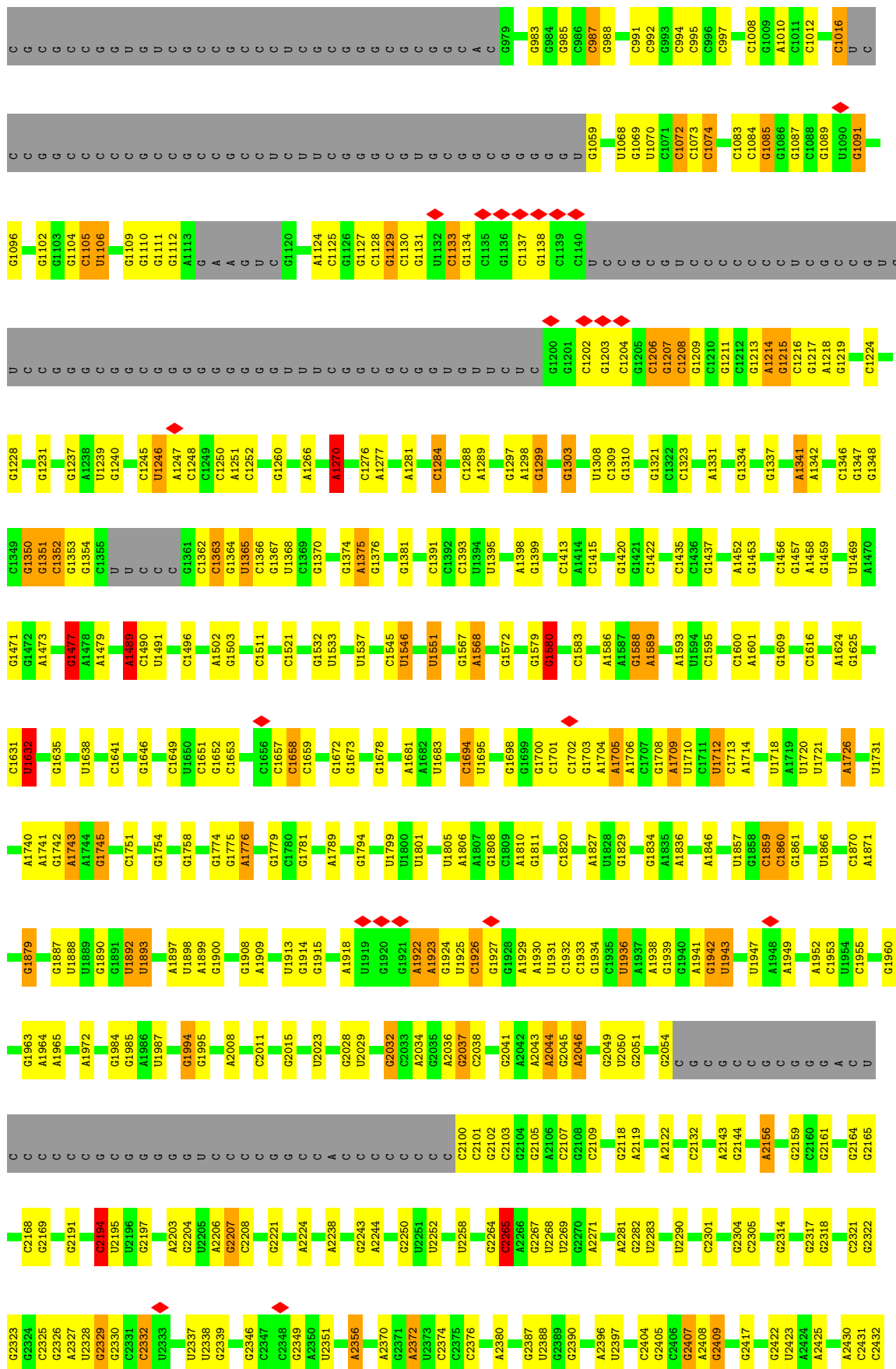




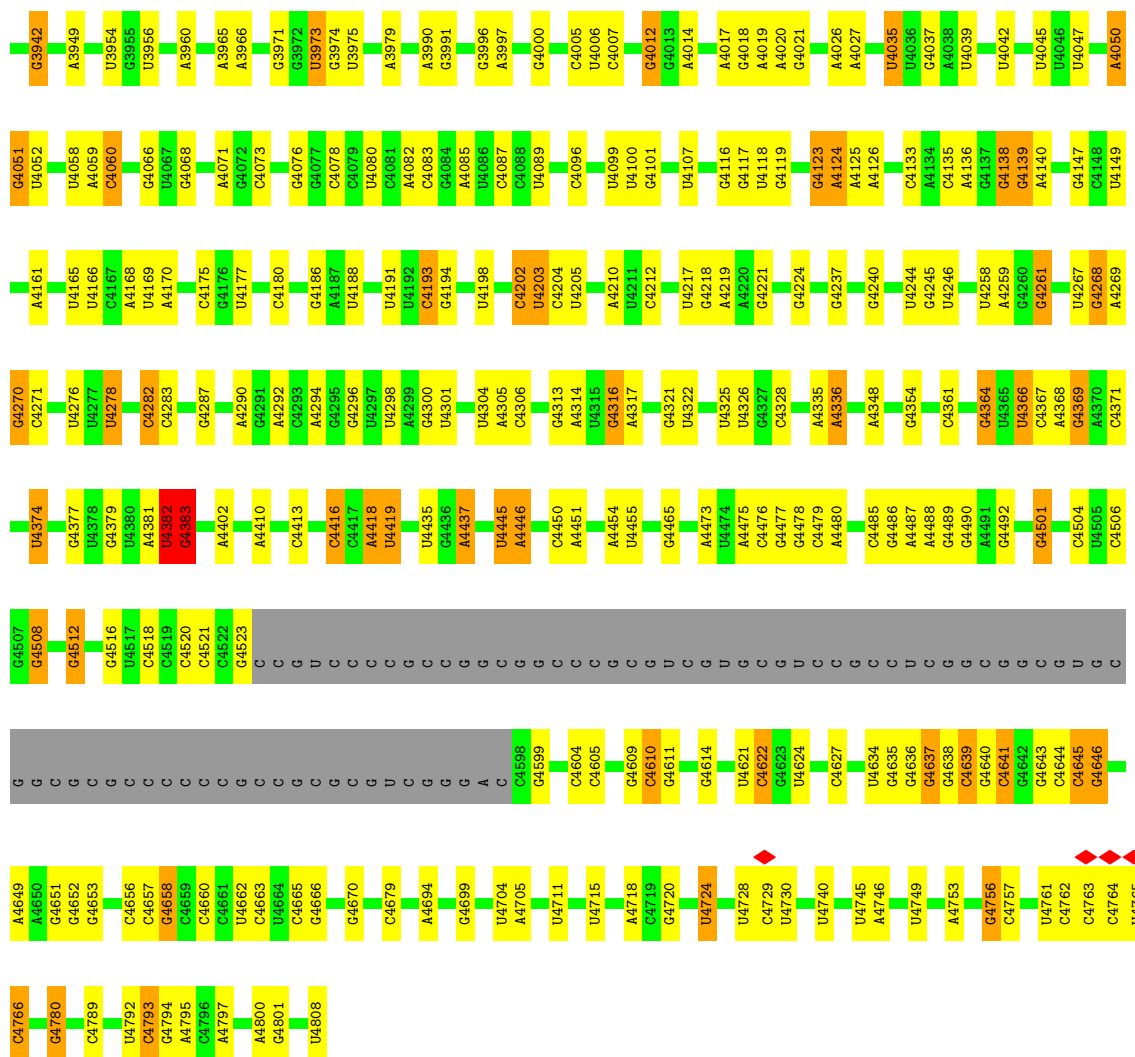
• Molecule 6: 28S rRNA

Chain L5:









• Molecule 7: 5S rRNA

Chain L7: 82% 17% .



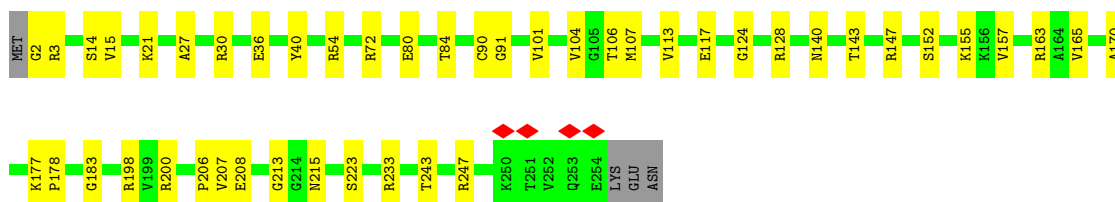
• Molecule 8: 5.8S rRNA

Chain L8: 78% 20% ..



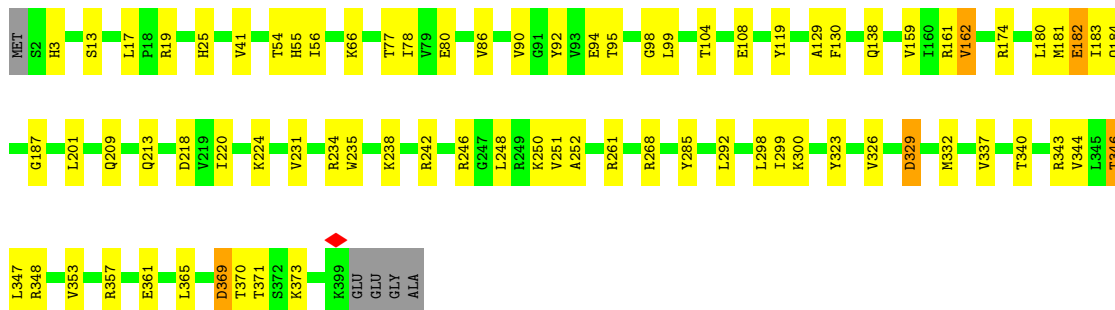
• Molecule 9: 60S ribosomal protein uL2

Chain LD: 81% 18% .



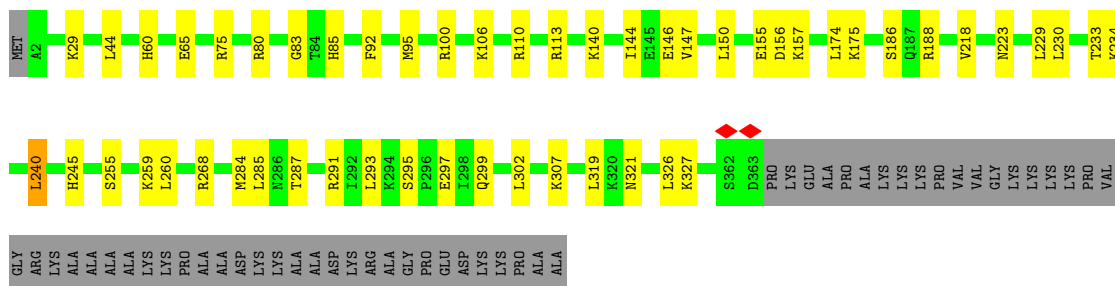
• Molecule 10: 60S ribosomal protein L3

Chain LE: 79% 18% ..



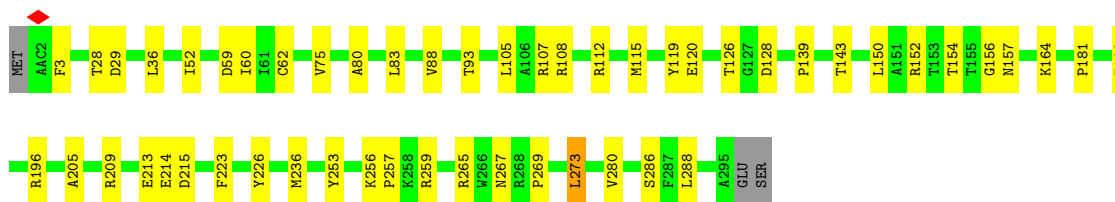
• Molecule 11: 60S ribosomal protein L4

Chain LF: 75% 12% 12%



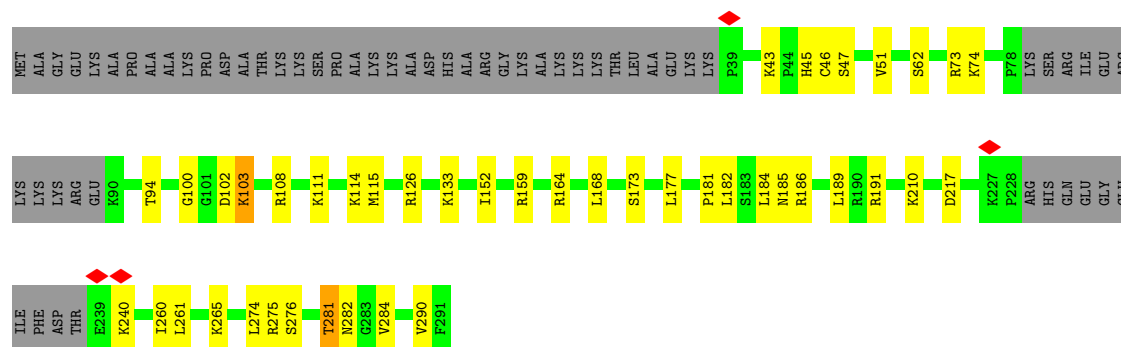
• Molecule 12: Ribosomal_L18_c domain-containing protein

Chain LG: 81% 17% .



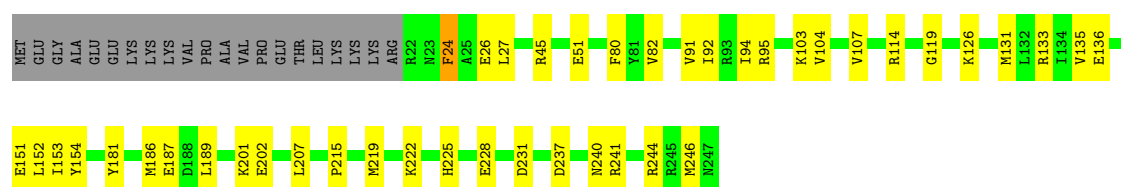
• Molecule 13: 60S ribosomal protein L6

Chain LH: 65% 14% 20%



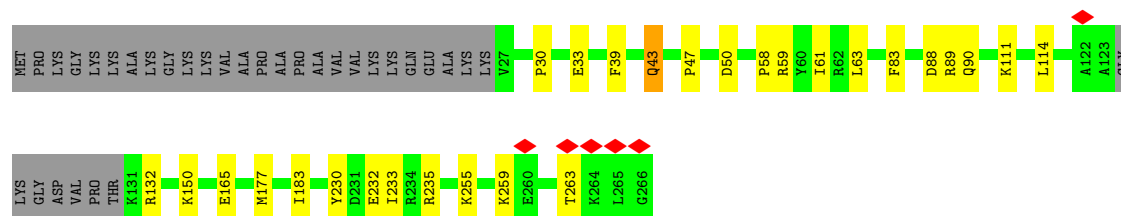
• Molecule 14: 60S ribosomal Protein uL30

Chain LI: 74% 17% 9%



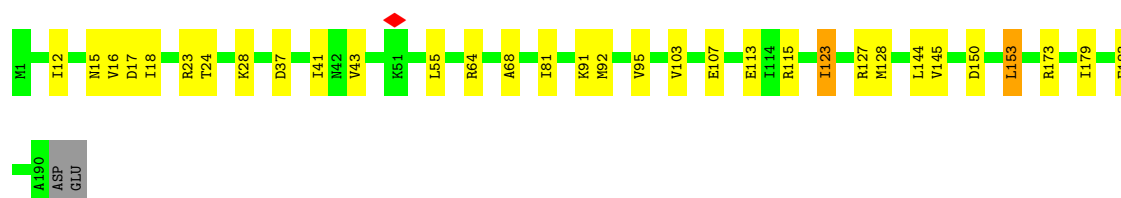
• Molecule 15: 60S ribosomal protein eL8

Chain LJ: 77% 10% 12%



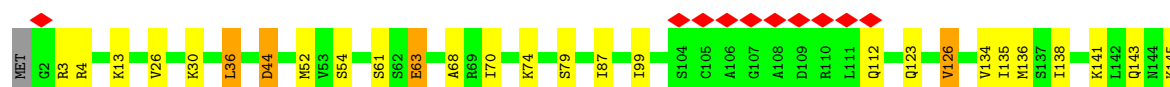
• Molecule 16: 60S ribosomal protein L9

Chain LK: 82% 16% ..



• Molecule 17: 60S ribosomal protein L10

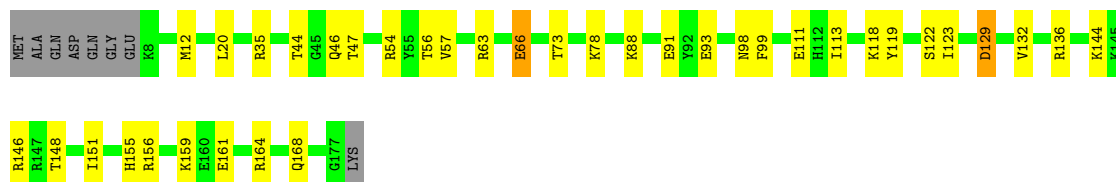
Chain LL: 5% 80% 17%





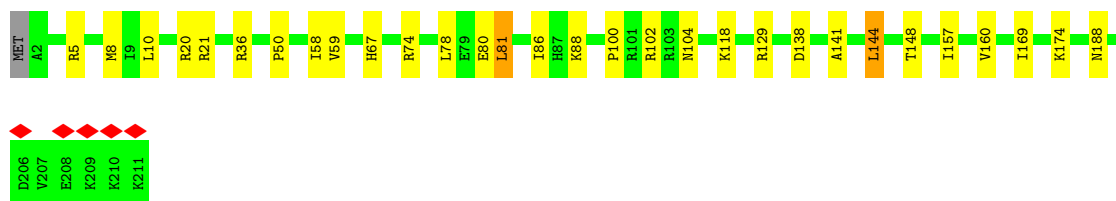
- Molecule 18: 60S ribosomal protein L11

Chain LM: 75% 20% . .



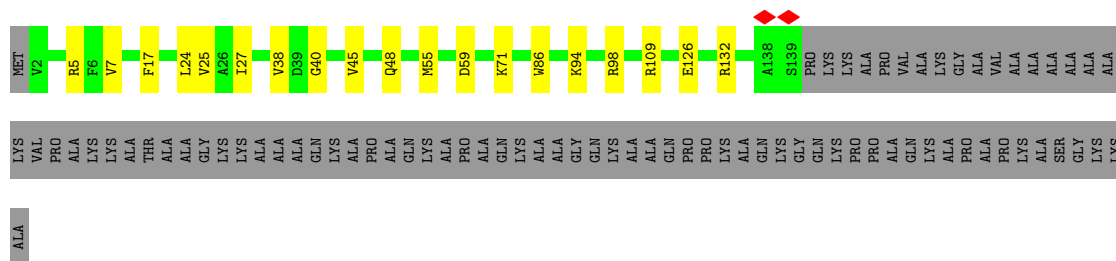
- Molecule 19: 60S ribosomal protein eL13

Chain LO: 85% 13% .



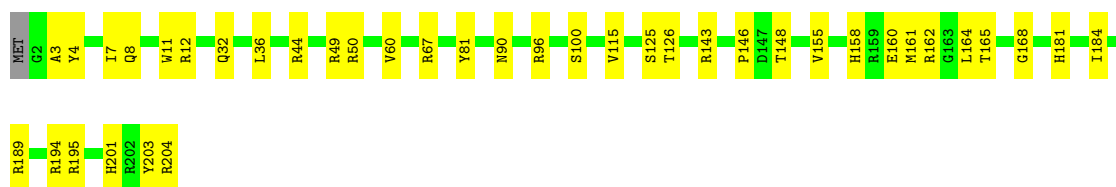
- Molecule 20: 60S ribosomal protein L14

Chain LP: 55% 9% 37%



- Molecule 21: 60S ribosomal protein L15

Chain LQ: 80% 19%



- Molecule 22: 60S ribosomal protein uL13

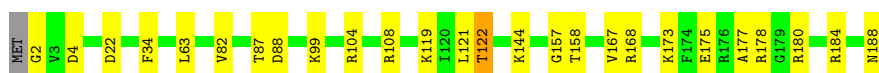
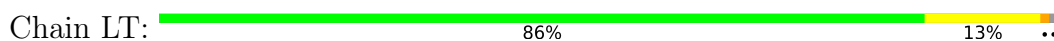
Chain LR: 87% 11% .



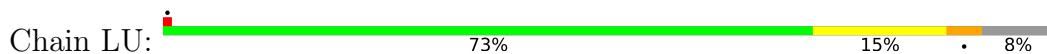
- Molecule 23: 60S ribosomal protein uL22



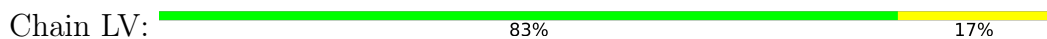
- Molecule 24: 60S ribosomal Protein eL18



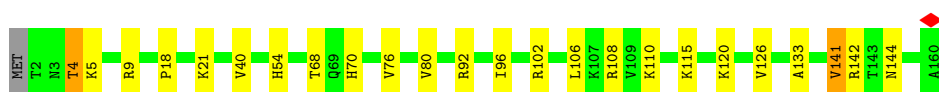
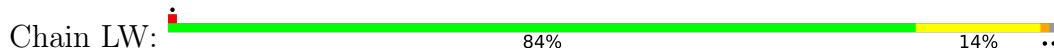
- Molecule 25: 60S ribosomal protein L19



- Molecule 26: 60S ribosomal protein eL20

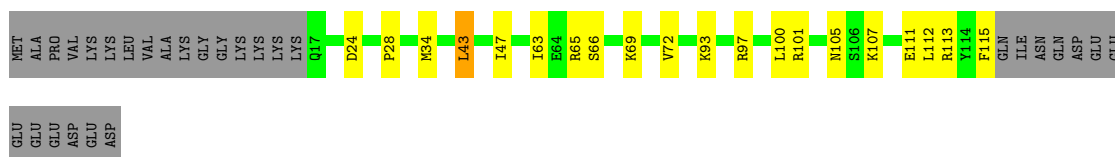


- Molecule 27: 60S ribosomal protein eL21



- Molecule 28: 60S ribosomal protein eL22

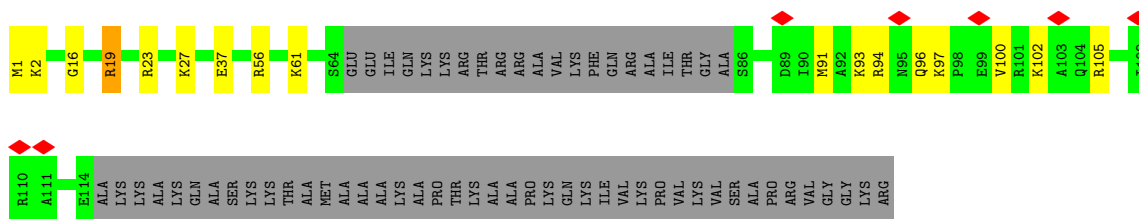




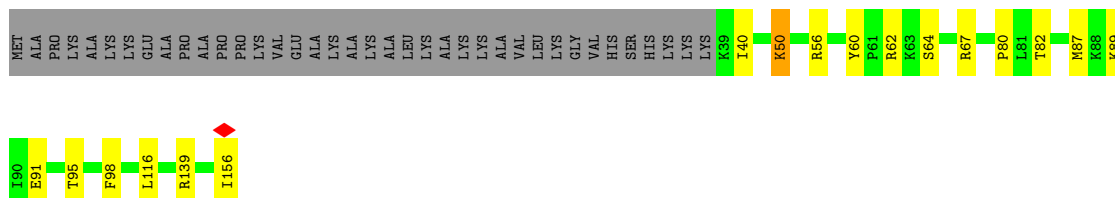
- Molecule 29: 60S ribosomal protein L23



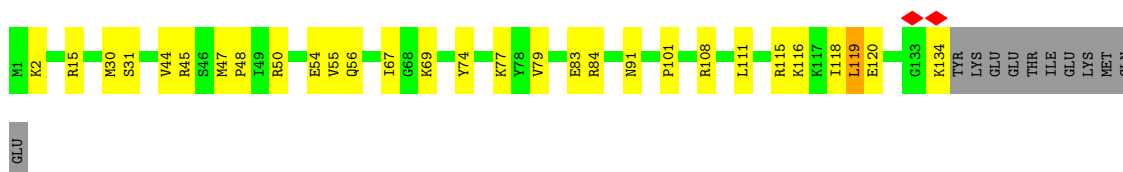
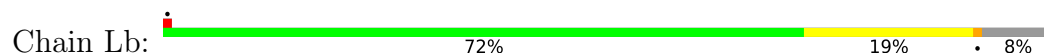
- Molecule 30: 60S ribosomal protein eL24



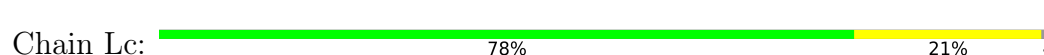
- Molecule 31: 60S ribosomal protein uL23



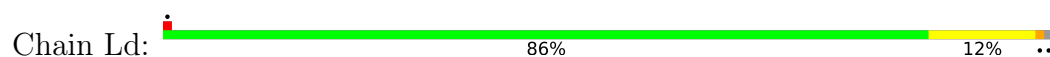
- Molecule 32: 60S ribosomal protein L26



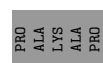
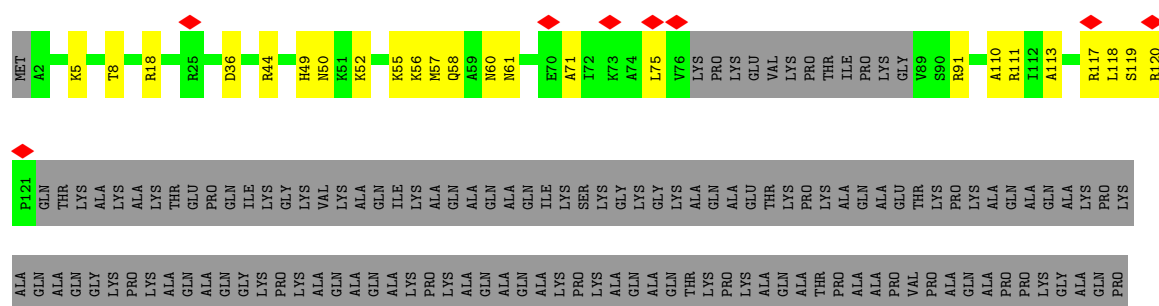
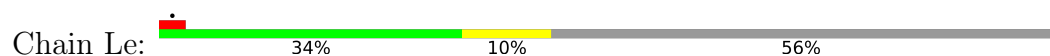
- Molecule 33: 60S ribosomal protein L27



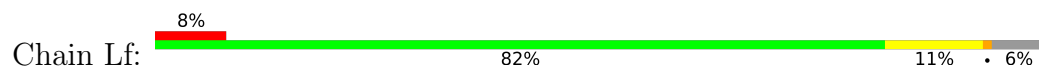
- Molecule 34: 60S ribosomal protein L27a



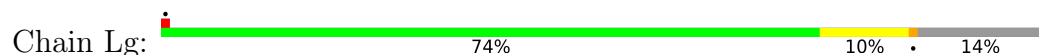
- Molecule 35: 60S ribosomal protein L29



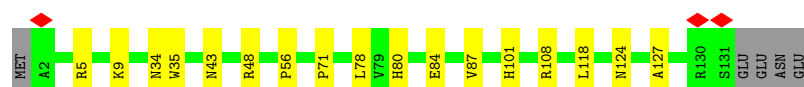
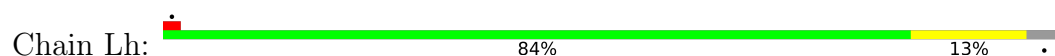
- Molecule 36: 60S ribosomal protein eL30



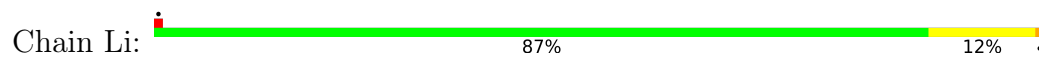
- Molecule 37: 60S ribosomal protein eL31



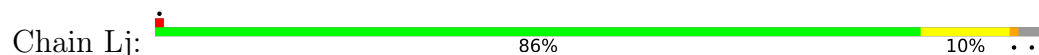
- Molecule 38: 60S ribosomal protein eL32



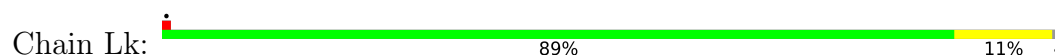
- Molecule 39: 60S ribosomal protein eL33



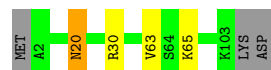
- Molecule 40: 60S ribosomal protein L34



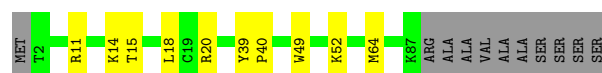
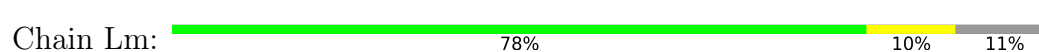
- Molecule 41: 60S ribosomal protein uL29



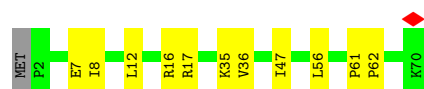
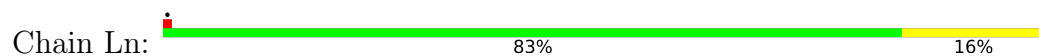
- Molecule 42: 60S ribosomal protein L36



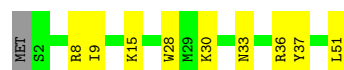
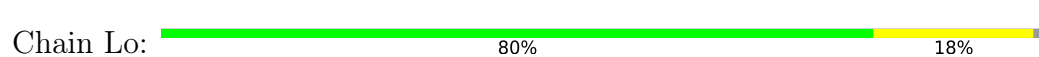
- Molecule 43: 60S ribosomal protein L37



- Molecule 44: 60S ribosomal protein eL38



- Molecule 45: 60S ribosomal protein eL39



- Molecule 46: 60S ribosomal protein L40

Chain Lp: 

MET GLN ILE PHE VAL SER THR LEU THR LYS THR ILE THR LEU GLU VAL PRO SER ASP THR ILE GLU ASN VAL LYS ALA LYS ILE GLN ASP LYS GLU ILE PRO PRO ASP GLN ARG LEU ILE PHE ALA GLY LYS GLN LEU GLU ASP GLY ARG THR LEU SER ASP TYR ASN

ILE GLN LYS GLU SER THR LEU HIS LEU VAL LYS ARG LEU ARG GLY I77 I78 L82 R83 N94 C99 Y100 P105 K114 K124 K125 K128

- Molecule 47: 60S ribosomal protein eL42

Chain Lq: 

MET Y2 F11 Y26 D31 Y34 A35 Q36 R39 Y48 K61 V67 E71 R78 R81 Q105 F106

- Molecule 48: 60S ribosomal protein eL43

Chain Lr: 

MET H2 G15 T16 R17 V28 I29 E30 T38 T45 V52 M61 Q92

- Molecule 49: 60S ribosomal protein eL28

Chain Ls: 

MET S2 L17 I18 K19 R20 S26 T27 E28 R30 D36 V61 V78 R79 T80 R87 R94 R98 K99 M108 I111 I117 S120 V126 K127 ARG LYS ARG ARG THR PRO ARG THR LYS SER

- Molecule 50: 60S ribosomal protein P0

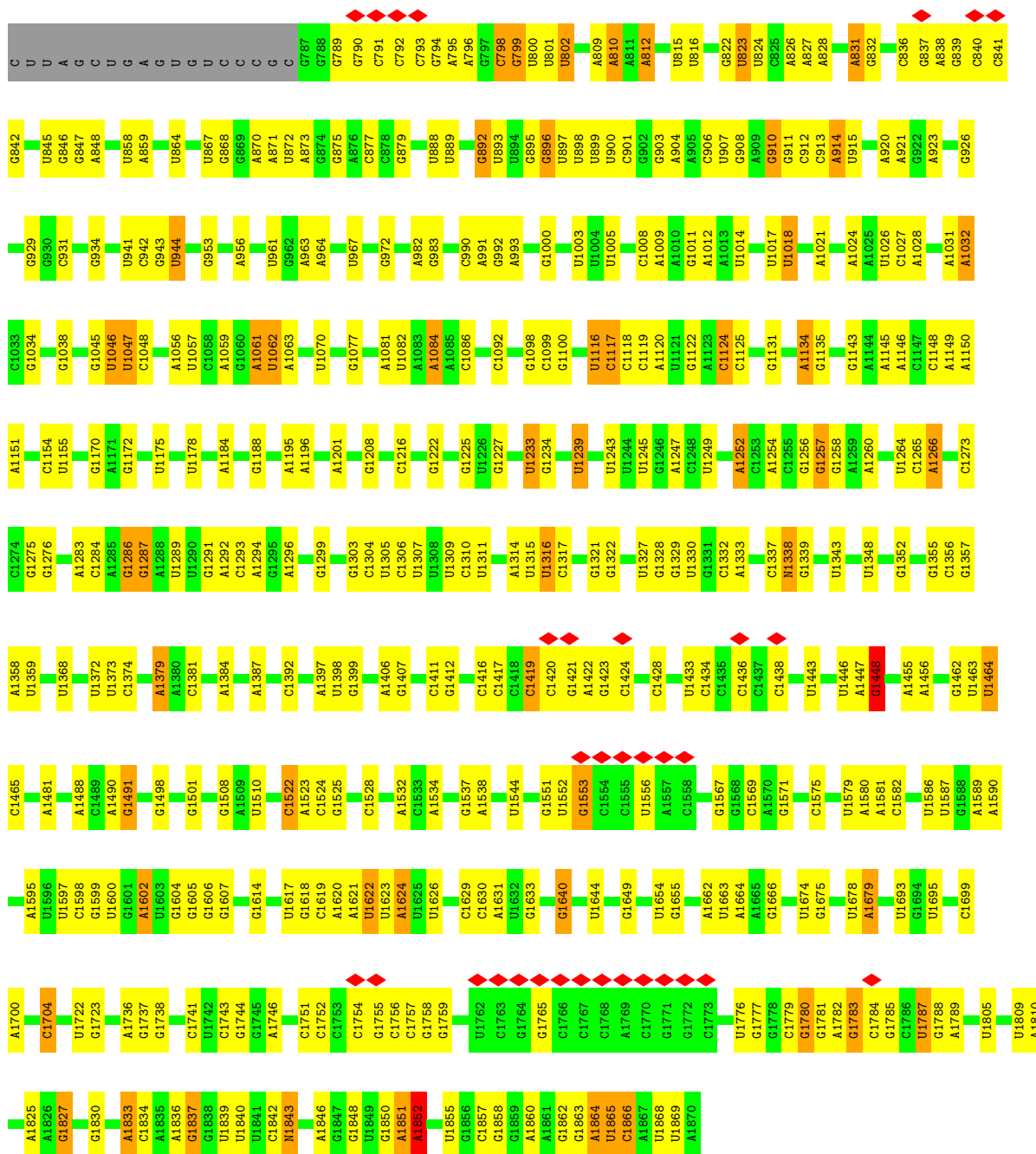
Chain Lt: 

MET PRO ARG GLU D5 Q19 L20 I29 V30 G31 A32 G36 S37 Q42 I43 R44 M45 S46 L54 R66 E70 F73 A74 L75 E76 K77 L78 L79 P80 H81 G84 G87 F88 W89 E93 D94 L95 T96 E97 I98 R99 D100 L103 A104 N105 LEU VAL PRO

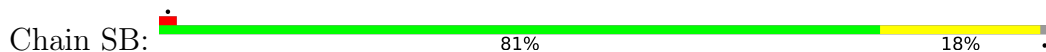
ALA ALA ARG ALA GLY ILE ILE PRO CYS VAL THR VAL PHE PHE ALA ASN THR GLN THR GLY LEU LEU GLY THR LYS THR THR THR LYS ILE SER ARG GLY THR ILE GLU ILE SER ASP VAL GLN LEU ILE LYS THR VAL THR GLY ASP LYS VAL CYS LEU GLN ILE

ALA SER GLU ALA ARG ALA LEU ASN MET LEU ASN ILE SER PRO PHE THR PHE PHE ALA GLY L187 L188 I189 Q190 Q191 V192 M195 G196 S197 I198 Y199 N200 PRO GLU ILE VAL ASP ILE THR ASP THR LEU HIS VAL ARG PHE LEU LEU GLY VAL VAL ALA ARG ASN VAL ALA SER THR ALA CYS LEU GLN ILE

GLY TYR PRO THR VAL SER VAL PRO HIS SER ILE ASN GLY TYR LYS ARG VAL LEU LEU ALA LEU SER PHE THR ALA VAL PRO



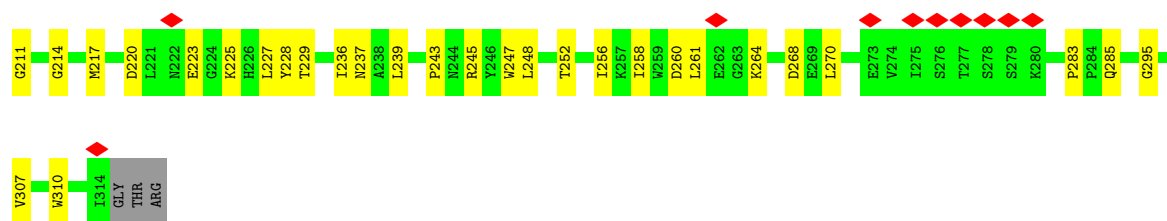
- Molecule 53: 40S ribosomal protein S27



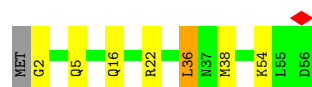
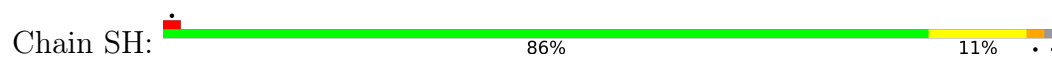
- Molecule 54: 40S ribosomal protein S28



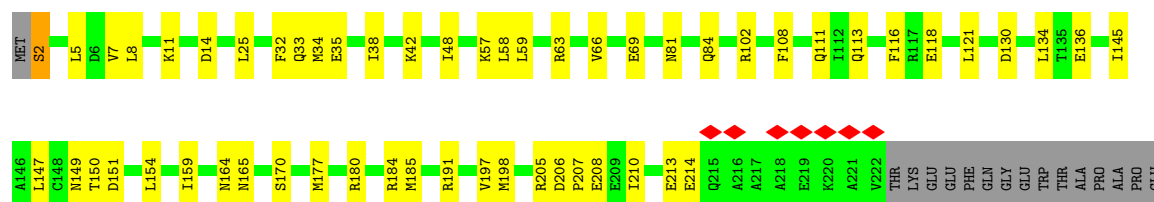




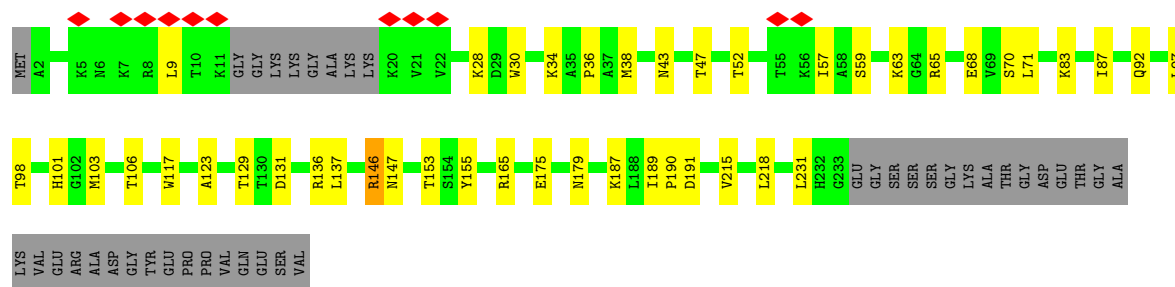
- Molecule 59: 40S ribosomal protein uS14

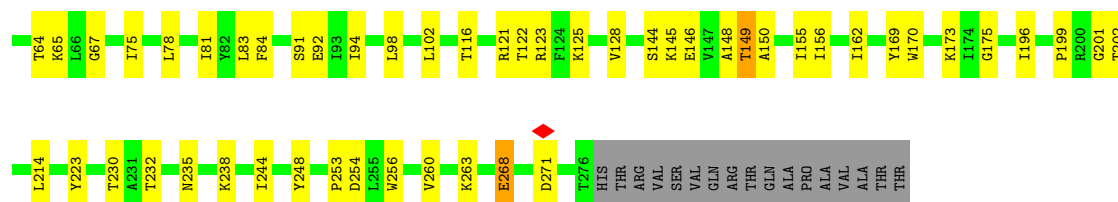


- Molecule 60: 40S ribosomal protein SA

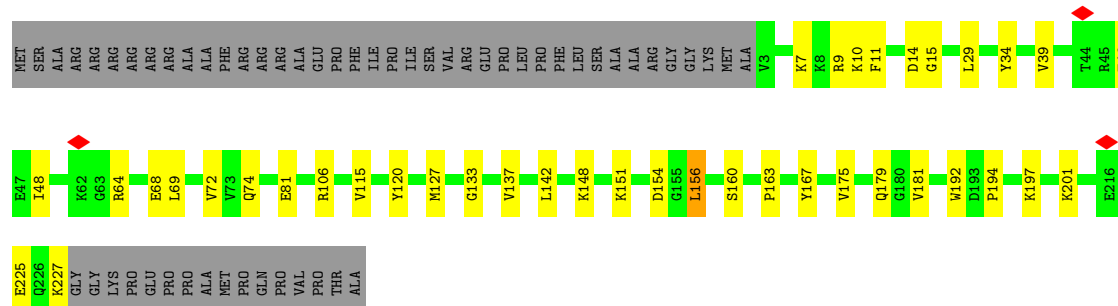


- Molecule 61: 40S ribosomal protein S3a

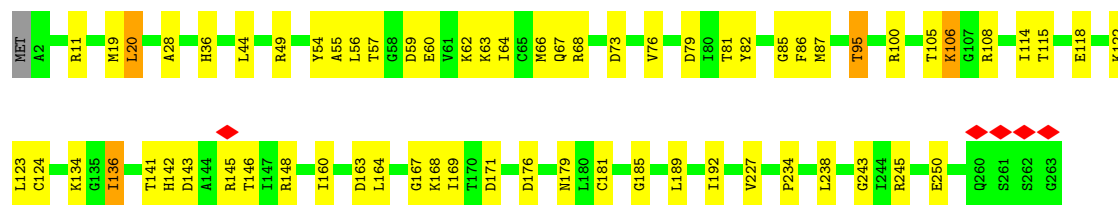




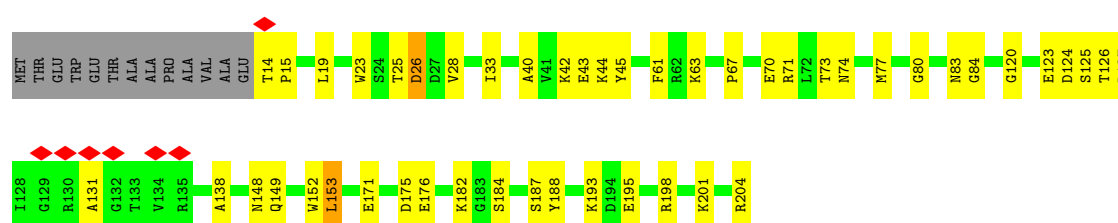
• Molecule 63: 40S ribosomal protein S3



• Molecule 64: 40S ribosomal protein eS4



• Molecule 65: Ribosomal protein S5

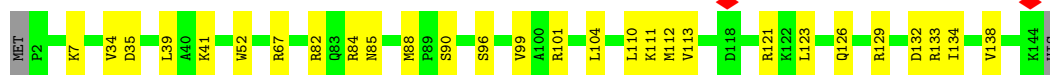


• Molecule 66: 40S ribosomal protein S6



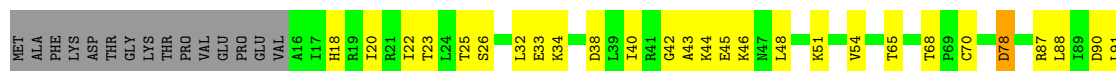


Chain Se:  79% 19%



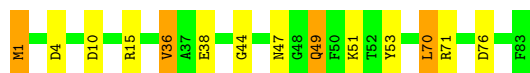
- Molecule 80: 40S ribosomal protein uS10

Chain Sf:  6% 58% 29% 13%



- Molecule 81: 40S ribosomal protein eS21

Chain Sg:  83% 12% 5%



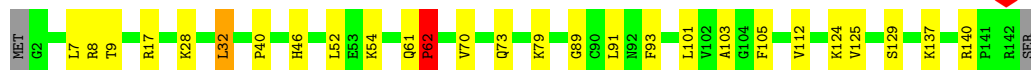
- Molecule 82: 40S ribosomal protein S15a

Chain Sh:  84% 15%



- Molecule 83: 40S ribosomal protein S23

Chain Si:  80% 17%

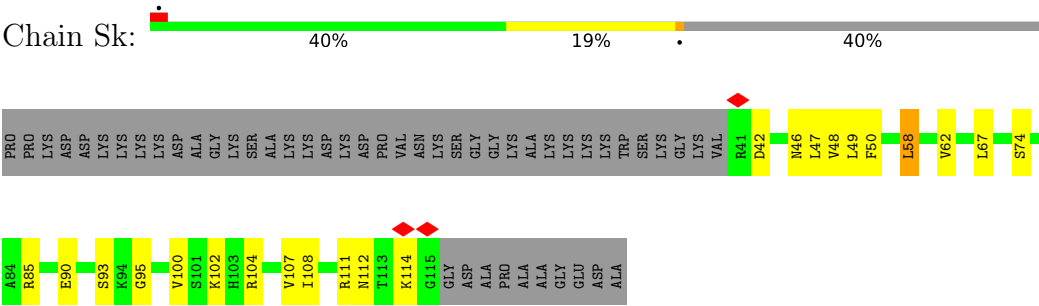


- Molecule 84: 40S ribosomal protein S24

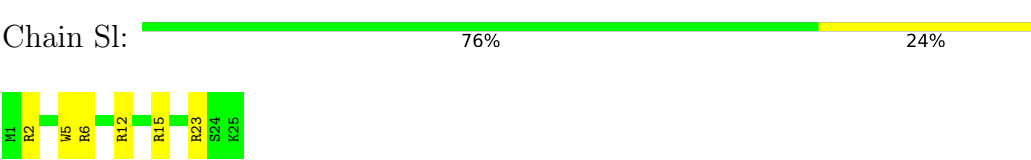
Chain Sj:  68% 28%



● Molecule 85: 40S ribosomal protein S25



● Molecule 86: 60S ribosomal protein L41



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	77142	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 ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.629	Depositor
Minimum map value	0.000	Depositor
Average map value	0.019	Depositor
Map value standard deviation	0.105	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	520.704, 520.704, 520.704	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.017, 1.017, 1.017	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 5MC, OMG, 7MG, OMC, PSU, NA, AME, SAC, GTP, MLZ, ZN, AAC, HIC, AYA, UR3, MA6, 1MA, NMM, 6MZ, OMU, HY3, A2M, 4AC, MG, K, M3L

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	E	0.10	0/3403	0.37	2/4656 (0.0%)
2	F	0.15	1/2136 (0.0%)	0.18	0/3325
3	I	0.10	0/4335	0.21	0/6738
4	S	0.09	0/1098	0.22	0/1710
5	B	0.15	0/1398	0.34	1/1905 (0.1%)
6	L5	0.18	1/87426 (0.0%)	0.19	0/136397
7	L7	0.16	0/2835	0.17	0/4418
8	L8	0.16	0/3635	0.17	0/5661
9	LD	0.16	0/1977	0.26	0/2651
10	LE	0.16	0/3261	0.26	0/4364
11	LF	0.16	0/2932	0.24	0/3939
12	LG	0.15	0/2437	0.23	0/3264
13	LH	0.13	0/1896	0.24	0/2541
14	LI	0.16	0/1922	0.24	0/2563
15	LJ	0.14	0/1908	0.23	0/2566
16	LK	0.14	0/1535	0.22	0/2063
17	LL	0.15	0/1756	0.25	0/2346
18	LM	0.13	0/1385	0.27	0/1852
19	LO	0.15	0/1733	0.24	0/2316
20	LP	0.15	0/1158	0.24	0/1547
21	LQ	0.18	0/1746	0.24	0/2338
22	LR	0.17	0/1662	0.24	0/2222
23	LS	0.16	0/1315	0.24	0/1763
24	LT	0.17	0/1539	0.27	0/2054
25	LU	0.13	0/1524	0.21	0/2013
26	LV	0.16	0/1497	0.23	0/2008
27	LW	0.15	0/1326	0.23	0/1770
28	LX	0.12	0/820	0.27	0/1100
29	LY	0.15	0/1048	0.25	0/1402
30	LZ	0.13	0/779	0.23	0/1034
31	La	0.14	0/984	0.24	0/1323

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	Lb	0.14	0/1132	0.22	0/1504
33	Lc	0.13	0/1130	0.20	0/1507
34	Ld	0.17	0/1191	0.26	0/1590
35	Le	0.13	0/884	0.28	0/1169
36	Lf	0.13	0/847	0.20	0/1134
37	Lg	0.15	0/903	0.23	0/1216
38	Lh	0.17	0/1088	0.24	0/1451
39	Li	0.17	0/903	0.28	0/1208
40	Lj	0.15	0/916	0.25	0/1220
41	Lk	0.13	0/1021	0.19	0/1348
42	Ll	0.12	0/841	0.21	0/1112
43	Lm	0.17	0/720	0.27	0/952
44	Ln	0.13	0/575	0.22	0/761
45	Lo	0.15	0/459	0.22	0/608
46	Lp	0.15	0/425	0.26	0/562
47	Lq	0.14	0/866	0.22	0/1141
48	Lr	0.14	0/718	0.24	0/953
49	Ls	0.16	0/1020	0.25	0/1366
50	Lt	0.09	0/565	0.26	0/783
51	Lx	0.09	0/969	0.26	0/1347
52	S2	0.16	0/40365	0.17	0/62915
53	SB	0.10	0/665	0.23	0/891
54	SC	0.10	0/497	0.22	0/666
55	SD	0.09	0/622	0.25	0/822
56	SE	0.08	0/462	0.24	0/607
57	SF	0.12	0/828	0.23	0/1109
58	SG	0.09	0/2493	0.25	0/3394
59	SH	0.12	0/470	0.23	0/623
60	SL	0.11	0/1771	0.23	0/2406
61	SM	0.11	0/1841	0.23	0/2459
62	SN	0.12	0/1742	0.24	0/2354
63	SO	0.11	0/1779	0.23	0/2395
64	SP	0.09	0/2118	0.24	0/2849
65	SQ	0.10	0/1531	0.25	0/2059
66	SR	0.09	0/1946	0.24	0/2590
67	SS	0.10	0/1552	0.24	0/2079
68	ST	0.10	0/1715	0.24	0/2287
69	SU	0.10	0/1550	0.25	0/2069
70	SV	0.11	0/834	0.25	0/1125
71	SW	0.11	0/1284	0.24	0/1717
72	SX	0.09	0/968	0.25	0/1296
73	SY	0.11	0/1232	0.22	0/1656
74	SZ	0.12	0/1029	0.27	0/1380

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
75	Sa	0.10	0/1069	0.23	0/1429
76	Sb	0.11	0/1142	0.29	0/1528
77	Sc	0.10	0/1094	0.24	0/1469
78	Sd	0.11	0/1226	0.26	0/1643
79	Se	0.09	0/1119	0.21	0/1498
80	Sf	0.11	0/831	0.24	0/1115
81	Sg	0.10	0/636	0.22	0/852
82	Sh	0.12	0/1051	0.23	0/1406
83	Si	0.12	0/1107	0.27	0/1475
84	Sj	0.09	0/1032	0.23	0/1371
85	Sk	0.09	0/604	0.25	0/810
86	Sl	0.14	0/240	0.16	0/305
All	All	0.16	2/242024 (0.0%)	0.21	3/355430 (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
83	Si	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	52	C	C4-N4	-5.20	1.23	1.33
6	L5	3517	A2M	O3'-P	5.04	1.61	1.56

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	571	PRO	N-CA-CB	11.87	110.78	102.35
1	E	546	PRO	N-CA-CB	7.34	110.95	103.25
5	B	704	LYS	CB-CA-C	-5.31	110.43	116.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
83	Si	62	HY3	Mainchain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3335	0	2684	42	0
2	F	1914	0	966	22	0
3	I	3882	0	1960	55	0
4	S	983	0	500	12	0
5	B	1379	0	1208	30	0
6	L5	80771	0	40889	609	0
7	L7	2538	0	1284	13	0
8	L8	3319	0	1684	12	0
9	LD	1939	0	2036	32	0
10	LE	3206	0	3353	51	0
11	LF	2886	0	3057	38	0
12	LG	2398	0	2430	35	0
13	LH	1859	0	2033	35	0
14	LI	1886	0	2008	29	0
15	LJ	1877	0	2023	18	0
16	LK	1516	0	1595	19	0
17	LL	1717	0	1762	27	0
18	LM	1362	0	1399	26	0
19	LO	1702	0	1820	22	0
20	LP	1137	0	1211	12	0
21	LQ	1701	0	1749	30	0
22	LR	1630	0	1776	18	0
23	LS	1288	0	1326	23	0
24	LT	1515	0	1634	22	0
25	LU	1508	0	1664	25	0
26	LV	1457	0	1492	17	0
27	LW	1298	0	1366	15	0
28	LX	806	0	827	14	0
29	LY	1034	0	1095	7	0
30	LZ	766	0	794	14	0
31	La	967	0	1040	13	0
32	Lb	1115	0	1205	21	0
33	Lc	1107	0	1182	22	0
34	Ld	1162	0	1205	15	0
35	Le	881	0	957	20	0
36	Lf	836	0	888	8	0
37	Lg	888	0	930	7	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
38	Lh	1070	0	1164	13	0
39	Li	884	0	924	8	0
40	Lj	906	0	998	10	0
41	Lk	1013	0	1147	9	0
42	Ll	830	0	916	3	0
43	Lm	705	0	737	7	0
44	Ln	569	0	637	9	0
45	Lo	447	0	480	7	0
46	Lp	431	0	471	7	0
47	Lq	863	0	929	8	0
48	Lr	708	0	756	8	0
49	Ls	1014	0	1083	11	0
50	Lt	567	0	251	5	0
51	Lx	972	0	429	5	0
52	S2	37825	0	19179	386	0
53	SB	651	0	672	10	0
54	SC	495	0	523	9	0
55	SD	610	0	634	29	0
56	SE	457	0	502	20	0
57	SF	814	0	862	21	0
58	SG	2436	0	2393	57	0
59	SH	459	0	448	7	0
60	SL	1743	0	1748	39	0
61	SM	1815	0	1908	29	0
62	SN	1706	0	1796	38	0
63	SO	1751	0	1846	26	0
64	SP	2076	0	2177	44	0
65	SQ	1509	0	1563	31	0
66	SR	1923	0	2089	54	0
67	SS	1529	0	1625	35	0
68	ST	1686	0	1772	49	0
69	SU	1525	0	1640	51	0
70	SV	810	0	836	13	0
71	SW	1262	0	1332	27	0
72	SX	958	0	993	31	0
73	SY	1208	0	1294	25	0
74	SZ	1016	0	1039	16	0
75	Sa	1048	0	1093	28	0
76	Sb	1124	0	1193	24	0
77	Sc	1080	0	1135	29	0
78	Sd	1217	0	1279	38	0
79	Se	1113	0	1145	20	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
80	Sf	821	0	883	22	0
81	Sg	640	0	633	14	0
82	Sh	1034	0	1078	16	0
83	Si	1099	0	1162	17	0
84	Sj	1015	0	1085	28	0
85	Sk	598	0	656	16	0
86	Sl	239	0	286	5	0
87	E	32	0	10	1	0
87	L7	32	0	11	3	0
88	F	6	0	4	0	0
89	I	2	0	0	0	0
89	L5	162	0	0	0	0
89	L7	2	0	0	0	0
89	L8	1	0	0	0	0
89	LD	1	0	0	0	0
89	LS	1	0	0	0	0
89	LY	1	0	0	0	0
89	S2	55	0	0	0	0
90	L5	17	0	0	0	0
90	S2	1	0	0	0	0
91	Lj	1	0	0	0	0
91	Lm	1	0	0	0	0
91	Lp	1	0	0	0	0
91	Lq	1	0	0	0	0
91	Lr	1	0	0	0	0
91	SD	1	0	0	0	0
91	SF	2	0	0	0	0
91	SH	1	0	0	0	0
92	SZ	1	0	0	0	0
93	L5	56	0	0	2	0
93	L7	1	0	0	0	0
93	L8	1	0	0	0	0
93	LD	1	0	0	0	0
93	LE	1	0	0	0	0
93	LH	1	0	0	0	0
93	LK	1	0	0	0	0
93	LL	1	0	0	0	0
93	LQ	1	0	0	0	0
93	LV	1	0	0	0	0
93	Ld	1	0	0	0	0
93	S2	34	0	0	0	0
93	SF	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
93	SH	1	0	0	0	0
93	SW	1	0	0	0	0
93	Sd	1	0	0	1	0
93	Se	1	0	0	0	0
93	Si	2	0	0	0	0
All	All	230266	0	168478	2304	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 6.

The worst 5 of 2304 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)
6:L5:278:G:H5''	21:LQ:8:GLN:HE22	1.45	0.80
76:Sb:102:GLU:HA	76:Sb:105:LYS:HB3	1.62	0.80
18:LM:164:ARG:HG2	18:LM:168:GLN:HE22	1.47	0.79
6:L5:1924:G:H1'	6:L5:1942:G:H5''	1.65	0.79
77:Sc:31:ASN:HD22	77:Sc:55:THR:HG22	1.47	0.78

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	E	509/616 (83%)	460 (90%)	46 (9%)	3 (1%)	21	51
5	B	200/854 (23%)	171 (86%)	28 (14%)	1 (0%)	24	55
9	LD	251/257 (98%)	241 (96%)	10 (4%)	0	100	100
10	LE	395/403 (98%)	386 (98%)	9 (2%)	0	100	100
11	LF	360/413 (87%)	355 (99%)	5 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
12	LG	291/297 (98%)	287 (99%)	4 (1%)	0	100	100
13	LH	226/291 (78%)	220 (97%)	6 (3%)	0	100	100
14	LI	224/247 (91%)	217 (97%)	7 (3%)	0	100	100
15	LJ	229/266 (86%)	227 (99%)	2 (1%)	0	100	100
16	LK	188/192 (98%)	187 (100%)	1 (0%)	0	100	100
17	LL	211/214 (99%)	209 (99%)	2 (1%)	0	100	100
18	LM	168/178 (94%)	162 (96%)	6 (4%)	0	100	100
19	LO	208/211 (99%)	204 (98%)	4 (2%)	0	100	100
20	LP	136/218 (62%)	135 (99%)	1 (1%)	0	100	100
21	LQ	201/204 (98%)	198 (98%)	3 (2%)	0	100	100
22	LR	197/203 (97%)	195 (99%)	2 (1%)	0	100	100
23	LS	157/184 (85%)	153 (98%)	4 (2%)	0	100	100
24	LT	185/188 (98%)	184 (100%)	1 (0%)	0	100	100
25	LU	178/196 (91%)	178 (100%)	0	0	100	100
26	LV	174/176 (99%)	171 (98%)	3 (2%)	0	100	100
27	LW	157/160 (98%)	156 (99%)	1 (1%)	0	100	100
28	LX	97/128 (76%)	94 (97%)	3 (3%)	0	100	100
29	LY	137/140 (98%)	134 (98%)	3 (2%)	0	100	100
30	LZ	89/157 (57%)	89 (100%)	0	0	100	100
31	La	116/156 (74%)	115 (99%)	1 (1%)	0	100	100
32	Lb	132/145 (91%)	129 (98%)	3 (2%)	0	100	100
33	Lc	133/136 (98%)	132 (99%)	1 (1%)	0	100	100
34	Ld	145/148 (98%)	140 (97%)	4 (3%)	1 (1%)	18	47
35	Le	103/245 (42%)	102 (99%)	1 (1%)	0	100	100
36	Lf	106/115 (92%)	106 (100%)	0	0	100	100
37	Lg	105/125 (84%)	105 (100%)	0	0	100	100
38	Lh	128/135 (95%)	127 (99%)	1 (1%)	0	100	100
39	Li	108/110 (98%)	108 (100%)	0	0	100	100
40	Lj	112/117 (96%)	112 (100%)	0	0	100	100
41	Lk	120/123 (98%)	118 (98%)	2 (2%)	0	100	100
42	Ll	100/105 (95%)	99 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
43	Lm	84/97 (87%)	83 (99%)	1 (1%)	0	100	100
44	Ln	67/70 (96%)	67 (100%)	0	0	100	100
45	Lo	48/51 (94%)	47 (98%)	1 (2%)	0	100	100
46	Lp	49/128 (38%)	49 (100%)	0	0	100	100
47	Lq	102/106 (96%)	101 (99%)	1 (1%)	0	100	100
48	Lr	89/92 (97%)	87 (98%)	2 (2%)	0	100	100
49	Ls	124/137 (90%)	121 (98%)	3 (2%)	0	100	100
50	Lt	111/318 (35%)	105 (95%)	5 (4%)	1 (1%)	14	41
51	Lx	190/217 (88%)	172 (90%)	17 (9%)	1 (0%)	24	55
53	SB	81/84 (96%)	80 (99%)	1 (1%)	0	100	100
54	SC	61/69 (88%)	60 (98%)	1 (2%)	0	100	100
55	SD	72/156 (46%)	71 (99%)	1 (1%)	0	100	100
56	SE	55/133 (41%)	54 (98%)	1 (2%)	0	100	100
57	SF	99/115 (86%)	99 (100%)	0	0	100	100
58	SG	311/317 (98%)	297 (96%)	14 (4%)	0	100	100
59	SH	53/56 (95%)	52 (98%)	1 (2%)	0	100	100
60	SL	219/295 (74%)	213 (97%)	6 (3%)	0	100	100
61	SM	220/264 (83%)	216 (98%)	4 (2%)	0	100	100
62	SN	218/293 (74%)	216 (99%)	2 (1%)	0	100	100
63	SO	223/281 (79%)	219 (98%)	4 (2%)	0	100	100
64	SP	260/263 (99%)	252 (97%)	8 (3%)	0	100	100
65	SQ	189/204 (93%)	183 (97%)	6 (3%)	0	100	100
66	SR	235/249 (94%)	232 (99%)	3 (1%)	0	100	100
67	SS	188/432 (44%)	183 (97%)	5 (3%)	0	100	100
68	ST	204/208 (98%)	197 (97%)	7 (3%)	0	100	100
69	SU	183/194 (94%)	176 (96%)	7 (4%)	0	100	100
70	SV	94/165 (57%)	91 (97%)	3 (3%)	0	100	100
71	SW	152/158 (96%)	148 (97%)	4 (3%)	0	100	100
72	SX	122/132 (92%)	116 (95%)	6 (5%)	0	100	100
73	SY	148/151 (98%)	146 (99%)	2 (1%)	0	100	100
74	SZ	134/151 (89%)	127 (95%)	6 (4%)	1 (1%)	18	47

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
75	Sa	126/145 (87%)	123 (98%)	3 (2%)	0	100	100
76	Sb	139/172 (81%)	132 (95%)	7 (5%)	0	100	100
77	Sc	132/135 (98%)	132 (100%)	0	0	100	100
78	Sd	146/152 (96%)	144 (99%)	2 (1%)	0	100	100
79	Se	140/145 (97%)	138 (99%)	2 (1%)	0	100	100
80	Sf	102/119 (86%)	99 (97%)	3 (3%)	0	100	100
81	Sg	81/83 (98%)	79 (98%)	2 (2%)	0	100	100
82	Sh	127/130 (98%)	126 (99%)	1 (1%)	0	100	100
83	Si	137/143 (96%)	136 (99%)	1 (1%)	0	100	100
84	Sj	123/130 (95%)	122 (99%)	1 (1%)	0	100	100
85	Sk	73/124 (59%)	72 (99%)	1 (1%)	0	100	100
86	Sl	23/25 (92%)	23 (100%)	0	0	100	100
All	All	12310/15142 (81%)	11992 (97%)	310 (2%)	8 (0%)	49	77

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	546	PRO
51	Lx	108	ASP
50	Lt	79	LEU
1	E	297	LEU
5	B	703	SER

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	E	227/524 (43%)	221 (97%)	6 (3%)	40	75
5	B	102/761 (13%)	99 (97%)	3 (3%)	37	73
9	LD	195/199 (98%)	191 (98%)	4 (2%)	47	79
10	LE	344/347 (99%)	328 (95%)	16 (5%)	23	57

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
11	LF	302/337 (90%)	297 (98%)	5 (2%)	53	83
12	LG	247/250 (99%)	242 (98%)	5 (2%)	48	80
13	LH	205/251 (82%)	201 (98%)	4 (2%)	48	80
14	LI	197/215 (92%)	191 (97%)	6 (3%)	36	72
15	LJ	199/223 (89%)	196 (98%)	3 (2%)	57	84
16	LK	169/171 (99%)	165 (98%)	4 (2%)	43	77
17	LL	180/181 (99%)	172 (96%)	8 (4%)	25	59
18	LM	143/149 (96%)	139 (97%)	4 (3%)	38	73
19	LO	175/176 (99%)	169 (97%)	6 (3%)	32	68
20	LP	117/161 (73%)	114 (97%)	3 (3%)	40	75
21	LQ	171/172 (99%)	168 (98%)	3 (2%)	51	82
22	LR	171/173 (99%)	171 (100%)	0	100	100
23	LS	139/163 (85%)	138 (99%)	1 (1%)	76	91
24	LT	164/165 (99%)	161 (98%)	3 (2%)	51	82
25	LU	159/175 (91%)	147 (92%)	12 (8%)	12	37
26	LV	154/154 (100%)	149 (97%)	5 (3%)	34	70
27	LW	139/140 (99%)	133 (96%)	6 (4%)	26	60
28	LX	88/113 (78%)	87 (99%)	1 (1%)	65	87
29	LY	106/107 (99%)	106 (100%)	0	100	100
30	LZ	79/126 (63%)	78 (99%)	1 (1%)	61	86
31	La	106/134 (79%)	104 (98%)	2 (2%)	50	81
32	Lb	124/135 (92%)	122 (98%)	2 (2%)	55	83
33	Lc	117/118 (99%)	116 (99%)	1 (1%)	70	89
34	Ld	119/120 (99%)	115 (97%)	4 (3%)	32	68
35	Le	87/183 (48%)	86 (99%)	1 (1%)	65	87
36	Lf	92/98 (94%)	90 (98%)	2 (2%)	45	78
37	Lg	98/110 (89%)	95 (97%)	3 (3%)	35	70
38	Lh	116/121 (96%)	113 (97%)	3 (3%)	40	75
39	Li	89/89 (100%)	86 (97%)	3 (3%)	32	68
40	Lj	98/100 (98%)	97 (99%)	1 (1%)	68	88
41	Lk	109/110 (99%)	106 (97%)	3 (3%)	38	73

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
42	Ll	86/89 (97%)	85 (99%)	1 (1%)	63	87
43	Lm	73/80 (91%)	70 (96%)	3 (4%)	27	62
44	Ln	64/65 (98%)	63 (98%)	1 (2%)	55	83
45	Lo	47/48 (98%)	47 (100%)	0	100	100
46	Lp	47/115 (41%)	47 (100%)	0	100	100
47	Lq	92/93 (99%)	91 (99%)	1 (1%)	65	87
48	Lr	74/75 (99%)	72 (97%)	2 (3%)	39	74
49	Ls	109/120 (91%)	104 (95%)	5 (5%)	24	58
53	SB	75/76 (99%)	73 (97%)	2 (3%)	39	74
54	SC	56/62 (90%)	53 (95%)	3 (5%)	20	51
55	SD	67/140 (48%)	65 (97%)	2 (3%)	36	72
56	SE	47/106 (44%)	46 (98%)	1 (2%)	47	79
57	SF	88/98 (90%)	87 (99%)	1 (1%)	65	87
58	SG	272/275 (99%)	263 (97%)	9 (3%)	33	69
59	SH	48/49 (98%)	47 (98%)	1 (2%)	47	79
60	SL	182/243 (75%)	178 (98%)	4 (2%)	45	78
61	SM	203/231 (88%)	199 (98%)	4 (2%)	48	80
62	SN	185/223 (83%)	180 (97%)	5 (3%)	39	74
63	SO	189/232 (82%)	184 (97%)	5 (3%)	40	75
64	SP	224/225 (100%)	216 (96%)	8 (4%)	31	66
65	SQ	161/170 (95%)	153 (95%)	8 (5%)	22	54
66	SR	207/218 (95%)	203 (98%)	4 (2%)	50	81
67	SS	170/360 (47%)	163 (96%)	7 (4%)	27	62
68	ST	178/180 (99%)	176 (99%)	2 (1%)	65	87
69	SU	161/168 (96%)	156 (97%)	5 (3%)	35	70
70	SV	87/136 (64%)	84 (97%)	3 (3%)	32	68
71	SW	139/142 (98%)	134 (96%)	5 (4%)	31	66
72	SX	104/108 (96%)	100 (96%)	4 (4%)	29	64
73	SY	130/131 (99%)	129 (99%)	1 (1%)	73	90
74	SZ	106/119 (89%)	101 (95%)	5 (5%)	23	57
75	Sa	114/130 (88%)	109 (96%)	5 (4%)	25	59

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
76	Sb	117/140 (84%)	113 (97%)	4 (3%)	32	68
77	Sc	120/121 (99%)	118 (98%)	2 (2%)	53	83
78	Sd	127/131 (97%)	125 (98%)	2 (2%)	55	83
79	Se	112/114 (98%)	110 (98%)	2 (2%)	51	82
80	Sf	94/107 (88%)	91 (97%)	3 (3%)	34	70
81	Sg	67/67 (100%)	64 (96%)	3 (4%)	24	58
82	Sh	112/113 (99%)	110 (98%)	2 (2%)	51	82
83	Si	112/114 (98%)	107 (96%)	5 (4%)	24	58
84	Sj	107/112 (96%)	107 (100%)	0	100	100
85	Sk	66/102 (65%)	64 (97%)	2 (3%)	36	72
86	Sl	24/24 (100%)	24 (100%)	0	100	100
All	All	10170/12403 (82%)	9904 (97%)	266 (3%)	41	75

5 of 266 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
74	SZ	40	THR
75	Sa	57	LEU
83	Si	32	LEU
26	LV	77	ASN
25	LU	158	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 78 such sidechains are listed below:

Mol	Chain	Res	Type
64	SP	197	ASN
76	Sb	80	GLN
66	SR	187	HIS
70	SV	28	HIS
79	Se	11	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	F	89/90 (98%)	26 (29%)	0
3	I	178/1072 (16%)	103 (57%)	1 (0%)

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	S	45/855 (5%)	22 (48%)	1 (2%)
52	S2	1764/1870 (94%)	267 (15%)	1 (0%)
6	L5	3753/4808 (78%)	524 (13%)	3 (0%)
7	L7	118/119 (99%)	8 (6%)	0
8	L8	155/158 (98%)	17 (10%)	0
All	All	6102/8972 (68%)	967 (15%)	6 (0%)

5 of 967 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	F	10	A
2	F	12	C
2	F	14	U
2	F	16	A
2	F	19	G

5 of 6 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
6	L5	1892	U
6	L5	4445	U
52	S2	1265	C
4	S	1111	A
3	I	66	G

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

218 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
6	A2M	L5	3557	6	22,25,26	1.72	7 (31%)	31,36,39	2.43	13 (41%)
6	OMU	L5	2680	6	19,22,23	0.99	2 (10%)	26,31,34	1.82	6 (23%)
52	OMC	S2	174	52	19,22,23	1.20	3 (15%)	26,31,34	0.88	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
52	PSU	S2	682	52	18,21,22	1.36	3 (16%)	22,30,33	1.86	5 (22%)
52	OMU	S2	121	52	19,22,23	0.99	2 (10%)	26,31,34	1.75	5 (19%)
6	PSU	L5	2351	6	18,21,22	1.34	3 (16%)	22,30,33	1.86	4 (18%)
6	OMC	L5	2194	6	19,22,23	1.11	2 (10%)	26,31,34	0.95	1 (3%)
52	OMU	S2	355	71,52	19,22,23	0.98	2 (10%)	26,31,34	1.83	5 (19%)
6	A2M	L5	398	6	22,25,26	1.71	7 (31%)	31,36,39	2.41	14 (45%)
6	OMC	L5	3433	6,90	19,22,23	1.16	2 (10%)	26,31,34	0.92	1 (3%)
52	OMG	S2	510	52,89	23,26,27	1.17	2 (8%)	33,38,41	1.89	10 (30%)
6	OMG	L5	1260	6	23,26,27	1.22	4 (17%)	33,38,41	1.92	10 (30%)
6	PSU	L5	1720	6	18,21,22	1.36	3 (16%)	22,30,33	1.82	5 (22%)
52	PSU	S2	823	52	18,21,22	1.40	3 (16%)	22,30,33	1.80	4 (18%)
6	PSU	L5	4246	6	18,21,22	1.39	3 (16%)	22,30,33	1.88	5 (22%)
6	PSU	L5	4740	6	18,21,22	1.38	3 (16%)	22,30,33	1.78	3 (13%)
6	PSU	L5	3490	6	18,21,22	1.35	3 (16%)	22,30,33	1.81	4 (18%)
6	PSU	L5	1638	6,90	18,21,22	1.34	3 (16%)	22,30,33	1.87	5 (22%)
52	PSU	S2	610	52	18,21,22	1.40	3 (16%)	22,30,33	1.79	4 (18%)
6	OMC	L5	2208	6,89	19,22,23	1.13	2 (10%)	26,31,34	0.91	1 (3%)
6	PSU	L5	3447	6	18,21,22	1.36	3 (16%)	22,30,33	1.81	4 (18%)
6	OMC	L5	3601	6	19,22,23	1.16	2 (10%)	26,31,34	0.89	0
6	PSU	L5	4435	6	18,21,22	1.35	3 (16%)	22,30,33	1.84	4 (18%)
52	A2M	S2	669	52,89	22,25,26	1.74	8 (36%)	31,36,39	2.61	15 (48%)
6	PSU	L5	1718	6	18,21,22	1.35	3 (16%)	22,30,33	1.82	4 (18%)
47	MLZ	Lq	53	47	8,9,10	0.46	0	4,9,11	0.08	0
6	PSU	L5	3652	6,89	18,21,22	1.31	3 (16%)	22,30,33	1.86	4 (18%)
52	PSU	S2	1005	52	18,21,22	1.63	5 (27%)	22,30,33	2.85	6 (27%)
52	PSU	S2	1178	52	18,21,22	1.39	3 (16%)	22,30,33	1.80	4 (18%)
6	PSU	L5	3371	6	18,21,22	1.33	3 (16%)	22,30,33	1.82	5 (22%)
52	OMG	S2	602	52	23,26,27	1.17	3 (13%)	33,38,41	1.86	10 (30%)
6	PSU	L5	3585	6,89	18,21,22	1.32	3 (16%)	22,30,33	1.83	4 (18%)
6	OMG	L5	4138	6	23,26,27	1.19	3 (13%)	33,38,41	1.89	10 (30%)
6	OMG	L5	2719	6	23,26,27	1.17	2 (8%)	33,38,41	2.04	12 (36%)
6	OMG	L5	4116	6	23,26,27	1.19	3 (13%)	33,38,41	1.87	10 (30%)
6	PSU	L5	4217	6	18,21,22	1.34	3 (16%)	22,30,33	1.81	5 (22%)
52	OMG	S2	1491	52,89	23,26,27	1.18	3 (13%)	33,38,41	1.88	9 (27%)
6	PSU	L5	3554	6	18,21,22	1.36	3 (16%)	22,30,33	1.81	5 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	A2M	L5	3562	6	22,25,26	1.71	7 (31%)	31,36,39	2.44	14 (45%)
6	A2M	L5	400	6	22,25,26	1.75	7 (31%)	31,36,39	2.51	14 (45%)
52	PSU	S2	1082	52	18,21,22	1.33	3 (16%)	22,30,33	1.82	4 (18%)
52	OMU	S2	1443	52,89	19,22,23	0.99	2 (10%)	26,31,34	1.74	4 (15%)
6	OMU	L5	4244	6	19,22,23	1.00	2 (10%)	26,31,34	1.83	5 (19%)
6	OMG	L5	4364	6	23,26,27	1.18	3 (13%)	33,38,41	1.88	10 (30%)
52	PSU	S2	407	52	18,21,22	1.39	3 (16%)	22,30,33	1.85	4 (18%)
6	OMG	L5	4383	6,90	23,26,27	1.20	3 (13%)	33,38,41	1.87	10 (30%)
52	OMG	S2	645	52	23,26,27	1.17	3 (13%)	33,38,41	1.92	10 (30%)
52	PSU	S2	1057	52	18,21,22	1.37	3 (16%)	22,30,33	1.84	5 (22%)
6	OMU	L5	2258	6	19,22,23	0.99	2 (10%)	26,31,34	1.77	5 (19%)
6	A2M	L5	1270	6	22,25,26	1.74	7 (31%)	31,36,39	2.56	16 (51%)
52	PSU	S2	119	52	18,21,22	1.41	3 (16%)	22,30,33	1.78	4 (18%)
52	PSU	S2	1239	52	18,21,22	1.37	3 (16%)	22,30,33	1.83	5 (22%)
11	AYA	LF	2	11	6,7,8	0.75	0	5,8,10	0.25	0
6	PSU	L5	4169	6	18,21,22	1.37	3 (16%)	22,30,33	1.85	5 (22%)
52	A2M	S2	591	52	22,25,26	1.72	6 (27%)	31,36,39	2.58	15 (48%)
35	MLZ	Le	5	35	8,9,10	0.47	0	4,9,11	0.13	0
52	PSU	S2	864	52	18,21,22	1.43	3 (16%)	22,30,33	1.81	4 (18%)
52	OMU	S2	172	52	19,22,23	0.99	2 (10%)	26,31,34	1.85	7 (26%)
52	OMC	S2	518	52	19,22,23	1.15	2 (10%)	26,31,34	0.91	1 (3%)
52	MA6	S2	1852	52	23,26,27	1.45	4 (17%)	34,38,41	2.33	13 (38%)
46	M3L	Lp	98	46	10,11,12	0.51	0	9,14,16	0.47	0
6	A2M	L5	3517	6	22,25,26	1.71	7 (31%)	31,36,39	2.62	15 (48%)
6	OMU	L5	4052	6	19,22,23	1.00	2 (10%)	26,31,34	1.74	4 (15%)
52	OMU	S2	429	52	19,22,23	1.02	2 (10%)	26,31,34	1.79	5 (19%)
6	OMC	L5	4202	6	19,22,23	1.14	2 (10%)	26,31,34	0.93	1 (3%)
8	PSU	L8	55	8	18,21,22	1.36	3 (16%)	22,30,33	1.84	4 (18%)
52	MA6	S2	1851	52	23,26,27	1.47	4 (17%)	34,38,41	2.33	13 (38%)
6	PSU	L5	3466	6	18,21,22	1.40	3 (16%)	22,30,33	1.86	4 (18%)
6	PSU	L5	1491	6	18,21,22	1.34	3 (16%)	22,30,33	1.85	4 (18%)
52	PSU	S2	816	52	18,21,22	1.43	3 (16%)	22,30,33	1.83	5 (22%)
6	PSU	L5	3583	6	18,21,22	1.33	3 (16%)	22,30,33	1.90	5 (22%)
6	PSU	L5	4039	6	18,21,22	1.35	3 (16%)	22,30,33	1.87	5 (22%)
6	PSU	L5	4045	6	18,21,22	1.35	3 (16%)	22,30,33	1.83	4 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	6MZ	L5	3966	6	22,25,26	1.39	4 (18%)	30,36,39	2.16	10 (33%)
6	OMC	L5	3619	6,22	19,22,23	1.14	2 (10%)	26,31,34	0.93	1 (3%)
6	PSU	L5	3500	6	18,21,22	1.36	3 (16%)	22,30,33	1.84	4 (18%)
52	OMU	S2	628	52	19,22,23	1.01	2 (10%)	26,31,34	1.82	6 (23%)
6	OMU	L5	4366	6	19,22,23	1.00	2 (10%)	26,31,34	1.81	6 (23%)
52	4AC	S2	1338	52	21,24,25	1.12	2 (9%)	29,34,37	1.21	3 (10%)
52	OMU	S2	1327	52,89	19,22,23	1.00	2 (10%)	26,31,34	1.79	5 (19%)
52	PSU	S2	218	52	18,21,22	1.38	3 (16%)	22,30,33	1.81	4 (18%)
6	OMG	L5	3524	6	23,26,27	1.19	3 (13%)	33,38,41	1.89	10 (30%)
6	PSU	L5	4099	6	18,21,22	1.35	3 (16%)	22,30,33	1.90	5 (22%)
6	OMG	L5	3942	6	23,26,27	1.18	3 (13%)	33,38,41	1.88	9 (27%)
52	OMG	S2	1448	52	23,26,27	1.18	3 (13%)	33,38,41	1.85	9 (27%)
6	PSU	L5	3502	6	18,21,22	1.39	3 (16%)	22,30,33	1.83	4 (18%)
52	A2M	S2	469	52	22,25,26	1.71	7 (31%)	31,36,39	2.47	15 (48%)
6	A2M	L5	2244	6	22,25,26	1.74	7 (31%)	31,36,39	2.51	15 (48%)
52	PSU	S2	36	52	18,21,22	1.41	3 (16%)	22,30,33	1.82	5 (22%)
6	PSU	L5	4382	6	18,21,22	1.42	4 (22%)	22,30,33	2.01	3 (13%)
6	A2M	L5	3450	6	22,25,26	1.72	7 (31%)	31,36,39	2.42	14 (45%)
6	OMC	L5	2667	6	19,22,23	1.14	2 (10%)	26,31,34	0.87	0
6	OMG	L5	3676	6	23,26,27	1.15	2 (8%)	33,38,41	1.87	9 (27%)
6	PSU	L5	4322	6	18,21,22	1.39	3 (16%)	22,30,33	1.79	5 (22%)
6	PSU	L5	4267	6,89	18,21,22	1.33	3 (16%)	22,30,33	1.93	4 (18%)
52	OMU	S2	1805	52	19,22,23	0.98	2 (10%)	26,31,34	1.80	5 (19%)
6	PSU	L5	3427	6	18,21,22	1.36	3 (16%)	22,30,33	1.81	5 (22%)
52	PSU	S2	34	52	18,21,22	1.39	3 (16%)	22,30,33	1.87	4 (18%)
52	6MZ	S2	1833	52,89	22,25,26	1.41	4 (18%)	30,36,39	2.17	9 (30%)
6	PSU	L5	3496	6	18,21,22	1.37	3 (16%)	22,30,33	1.84	5 (22%)
6	OMG	L5	4369	6	23,26,27	1.19	2 (8%)	33,38,41	1.87	9 (27%)
6	A2M	L5	3599	6	22,25,26	1.75	7 (31%)	31,36,39	2.52	16 (51%)
6	OMU	L5	3973	6	19,22,23	1.00	2 (10%)	26,31,34	1.83	6 (23%)
52	PSU	S2	802	52	18,21,22	1.40	3 (16%)	22,30,33	1.77	4 (18%)
6	PSU	L5	4058	6	18,21,22	1.37	3 (16%)	22,30,33	1.80	5 (22%)
52	A2M	S2	159	52	22,25,26	1.70	6 (27%)	31,36,39	2.40	14 (45%)
52	PSU	S2	1245	52	18,21,22	1.36	3 (16%)	22,30,33	1.85	4 (18%)
81	AME	Sg	1	81	9,10,11	0.47	0	9,11,13	0.88	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
52	PSU	S2	687	52	18,21,22	1.41	3 (16%)	22,30,33	1.84	5 (22%)
6	OMC	L5	2265	6,89	19,22,23	1.12	2 (10%)	26,31,34	0.90	2 (7%)
6	PSU	L5	1537	6	18,21,22	1.38	3 (16%)	22,30,33	1.81	4 (18%)
6	OMC	L5	2647	6	19,22,23	1.14	2 (10%)	26,31,34	0.88	1 (3%)
6	PSU	L5	3462	6	18,21,22	1.38	3 (16%)	22,30,33	1.80	5 (22%)
6	PSU	L5	3369	6,90,89	18,21,22	1.29	3 (16%)	22,30,33	1.87	4 (18%)
52	A2M	S2	577	52	22,25,26	1.71	7 (31%)	31,36,39	2.46	16 (51%)
83	HY3	Si	62	83	6,8,9	1.32	1 (16%)	5,10,12	1.43	1 (20%)
6	OMG	L5	3631	6	23,26,27	1.21	4 (17%)	33,38,41	1.90	10 (30%)
6	PSU	L5	1632	6	18,21,22	1.41	3 (16%)	22,30,33	1.74	4 (18%)
52	PSU	S2	1047	52	18,21,22	1.39	3 (16%)	22,30,33	1.79	4 (18%)
6	PSU	L5	1731	6	18,21,22	1.35	3 (16%)	22,30,33	1.83	4 (18%)
6	PSU	L5	3576	6	18,21,22	1.36	3 (16%)	22,30,33	1.77	4 (18%)
6	A2M	L5	4336	6	22,25,26	1.74	7 (31%)	31,36,39	2.51	15 (48%)
52	A2M	S2	1384	52	22,25,26	1.74	7 (31%)	31,36,39	2.46	15 (48%)
6	PSU	L5	3494	6	18,21,22	1.39	3 (16%)	22,30,33	1.78	5 (22%)
10	HIC	LE	245	10	10,11,12	0.55	0	8,14,16	0.48	0
52	PSU	S2	1368	52	18,21,22	1.41	3 (16%)	22,30,33	1.85	5 (22%)
52	PSU	S2	1046	52	18,21,22	1.35	3 (16%)	22,30,33	1.88	4 (18%)
6	PSU	L5	4374	6	18,21,22	1.36	3 (16%)	22,30,33	1.82	5 (22%)
52	PSU	S2	1644	52,89	18,21,22	1.36	3 (16%)	22,30,33	1.83	4 (18%)
52	PSU	S2	93	52	18,21,22	1.42	3 (16%)	22,30,33	1.76	3 (13%)
52	PSU	S2	109	52	18,21,22	1.42	3 (16%)	22,30,33	1.80	5 (22%)
52	A2M	S2	485	52	22,25,26	1.73	7 (31%)	31,36,39	2.54	15 (48%)
6	A2M	L5	3492	6,89	22,25,26	1.74	7 (31%)	31,36,39	2.47	14 (45%)
52	PSU	S2	1693	52	18,21,22	1.36	3 (16%)	22,30,33	1.86	5 (22%)
6	OMU	L5	3657	6	19,22,23	1.01	2 (10%)	26,31,34	1.82	5 (19%)
6	5MC	L5	4193	6	18,22,23	1.24	3 (16%)	26,32,35	1.53	3 (11%)
52	OMC	S2	1392	52	19,22,23	1.16	2 (10%)	26,31,34	0.94	1 (3%)
52	PSU	S2	652	52	18,21,22	1.39	3 (16%)	22,30,33	1.84	4 (18%)
52	PSU	S2	1348	52	18,21,22	1.38	3 (16%)	22,30,33	1.81	3 (13%)
52	OMU	S2	116	52	19,22,23	1.01	2 (10%)	26,31,34	1.80	6 (23%)
52	PSU	S2	1233	52	18,21,22	1.37	3 (16%)	22,30,33	1.81	5 (22%)
52	PSU	S2	573	52	18,21,22	1.43	3 (16%)	22,30,33	1.84	4 (18%)
6	PSU	L5	4166	6	18,21,22	1.38	3 (16%)	22,30,33	1.79	4 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	PSU	L5	4177	6	18,21,22	1.33	3 (16%)	22,30,33	1.86	5 (22%)
52	OMC	S2	1704	52,89	19,22,23	1.13	2 (10%)	26,31,34	0.89	1 (3%)
6	OMG	L5	3476	6	23,26,27	1.19	3 (13%)	33,38,41	1.87	9 (27%)
6	OMG	L5	4240	6	23,26,27	1.20	3 (13%)	33,38,41	1.87	10 (30%)
52	A2M	S2	27	52	22,25,26	1.74	7 (31%)	31,36,39	2.46	14 (45%)
6	UR3	L5	4276	6	19,22,23	1.30	3 (15%)	26,32,35	1.20	2 (7%)
6	OMC	L5	3540	6	19,22,23	1.10	2 (10%)	26,31,34	0.89	2 (7%)
52	OMG	S2	684	52	23,26,27	1.18	3 (13%)	33,38,41	1.90	10 (30%)
52	OMG	S2	868	52	23,26,27	1.16	2 (8%)	33,38,41	1.87	9 (27%)
6	OMG	L5	1477	6	23,26,27	1.21	3 (13%)	33,38,41	1.95	10 (30%)
6	PSU	L5	4203	6	18,21,22	1.33	3 (16%)	22,30,33	1.81	4 (18%)
52	PSU	S2	105	52	18,21,22	1.39	3 (16%)	22,30,33	1.81	4 (18%)
6	OMC	L5	1820	6,89	19,22,23	1.15	2 (10%)	26,31,34	0.89	0
6	A2M	L5	3456	6	22,25,26	1.71	7 (31%)	31,36,39	2.41	13 (41%)
52	A2M	S2	1679	52	22,25,26	1.71	7 (31%)	31,36,39	2.43	13 (41%)
6	PSU	L5	1799	6	18,21,22	1.38	3 (16%)	22,30,33	1.86	5 (22%)
6	PSU	L5	1801	6	18,21,22	1.35	3 (16%)	22,30,33	1.83	3 (13%)
6	A2M	L5	2658	6,89	22,25,26	1.74	7 (31%)	31,36,39	2.49	15 (48%)
6	PSU	L5	4149	6	18,21,22	1.33	3 (16%)	22,30,33	1.93	5 (22%)
52	PSU	S2	1446	52	18,21,22	1.40	3 (16%)	22,30,33	1.85	4 (18%)
6	PSU	L5	1721	6	18,21,22	1.35	3 (16%)	22,30,33	1.83	5 (22%)
6	PSU	L5	4188	6	18,21,22	1.33	3 (16%)	22,30,33	1.93	4 (18%)
6	PSU	L5	4107	6	18,21,22	1.33	3 (16%)	22,30,33	1.79	3 (13%)
6	OMC	L5	2704	6	19,22,23	1.14	2 (10%)	26,31,34	0.86	0
6	PSU	L5	4711	6	18,21,22	1.35	3 (16%)	22,30,33	1.82	5 (22%)
6	1MA	L5	1266	6,89	21,25,26	1.82	5 (23%)	31,37,40	1.92	5 (16%)
6	A2M	L5	1810	6,89	22,25,26	1.76	7 (31%)	31,36,39	2.52	15 (48%)
6	OMG	L5	3359	6	23,26,27	1.19	3 (13%)	33,38,41	1.89	10 (30%)
6	PSU	L5	4749	6	18,21,22	1.39	3 (16%)	22,30,33	1.84	4 (18%)
6	A2M	L5	2630	6	22,25,26	1.75	7 (31%)	31,36,39	2.63	17 (54%)
52	A2M	S2	513	52	22,25,26	1.72	7 (31%)	31,36,39	2.43	13 (41%)
52	PSU	S2	650	52	18,21,22	1.38	3 (16%)	22,30,33	1.83	4 (18%)
52	PSU	S2	1626	52	18,21,22	1.40	3 (16%)	22,30,33	1.79	5 (22%)
6	PSU	L5	4419	6	18,21,22	1.34	3 (16%)	22,30,33	1.84	5 (22%)
52	A2M	S2	1032	52	22,25,26	1.71	7 (31%)	31,36,39	2.46	14 (45%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	PSU	L8	69	8	18,21,22	1.36	3 (16%)	22,30,33	1.90	5 (22%)
6	A2M	L5	4317	6	22,25,26	1.75	7 (31%)	31,36,39	2.46	14 (45%)
52	OMC	S2	463	52	19,22,23	1.16	2 (10%)	26,31,34	0.90	0
6	OMC	L5	3573	6	19,22,23	1.13	2 (10%)	26,31,34	0.89	2 (7%)
52	OMU	S2	1289	52,55	19,22,23	0.97	1 (5%)	26,31,34	1.76	5 (19%)
52	A2M	S2	99	52,89	22,25,26	1.71	6 (27%)	31,36,39	2.40	13 (41%)
6	5MC	L5	3514	6,89	18,22,23	1.27	3 (16%)	26,32,35	1.26	4 (15%)
6	OMG	L5	4245	6	23,26,27	1.20	4 (17%)	33,38,41	1.87	10 (30%)
52	PSU	S2	815	52	18,21,22	1.38	3 (16%)	22,30,33	1.83	4 (18%)
6	PSU	L5	4325	6	18,21,22	1.35	3 (16%)	22,30,33	1.85	5 (22%)
52	7MG	S2	1640	52	22,26,27	1.22	2 (9%)	29,39,42	2.14	9 (31%)
6	OMG	L5	2267	6	23,26,27	1.19	2 (8%)	33,38,41	1.85	9 (27%)
52	PSU	S2	967	52	18,21,22	1.40	3 (16%)	22,30,33	1.79	4 (18%)
52	A2M	S2	166	52	22,25,26	1.71	6 (27%)	31,36,39	2.46	15 (48%)
6	PSU	L5	1683	6,90	18,21,22	1.36	3 (16%)	22,30,33	1.82	4 (18%)
6	OMG	L5	3974	6	23,26,27	1.19	3 (13%)	33,38,41	1.86	9 (27%)
6	PSU	L5	2475	6	18,21,22	1.37	3 (16%)	22,30,33	1.82	5 (22%)
6	OMC	L5	4282	6	19,22,23	1.15	2 (10%)	26,31,34	0.91	1 (3%)
6	A2M	L5	4269	6,89	22,25,26	1.74	7 (31%)	31,36,39	2.45	14 (45%)
52	4AC	S2	1843	52,86	21,24,25	1.15	2 (9%)	29,34,37	1.10	2 (6%)
8	OMG	L8	75	8	23,26,27	1.19	3 (13%)	33,38,41	1.85	9 (27%)
52	PSU	S2	867	52	18,21,22	1.44	3 (16%)	22,30,33	1.81	4 (18%)
6	A2M	L5	1479	6	22,25,26	1.72	7 (31%)	31,36,39	2.54	14 (45%)
6	OMG	L5	2207	6	23,26,27	1.18	3 (13%)	33,38,41	2.01	10 (30%)
6	PSU	L5	4278	6	18,21,22	1.38	3 (16%)	22,30,33	1.79	4 (18%)
52	PSU	S2	210	52	18,21,22	1.45	3 (16%)	22,30,33	1.84	4 (18%)
49	SAC	Ls	2	49	7,8,9	0.53	0	8,9,11	0.84	1 (12%)
6	PSU	L5	4298	6	18,21,22	1.34	3 (16%)	22,30,33	1.83	4 (18%)
78	SAC	Sd	2	78	7,8,9	0.52	0	8,9,11	0.90	1 (12%)
52	PSU	S2	1175	52	18,21,22	1.36	3 (16%)	22,30,33	1.84	5 (22%)
52	OMG	S2	1329	52	23,26,27	1.17	2 (8%)	33,38,41	1.86	10 (30%)
79	NMM	Se	67	79	9,11,12	1.56	1 (11%)	6,12,14	3.48	2 (33%)
6	A2M	L5	2206	6,89	22,25,26	1.76	7 (31%)	31,36,39	2.47	15 (48%)
6	OMC	L5	1284	6	19,22,23	1.16	2 (10%)	26,31,34	0.86	0
6	PSU	L5	4042	6	18,21,22	1.36	3 (16%)	22,30,33	1.87	4 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
52	OMG	S2	437	52	23,26,27	1.16	2 (8%)	33,38,41	1.88	9 (27%)
60	SAC	SL	2	60	7,8,9	0.53	0	8,9,11	0.86	1 (12%)
6	A2M	L5	1489	6,89	22,25,26	1.75	7 (31%)	31,36,39	2.42	12 (38%)
6	PSU	L5	3616	6	18,21,22	1.34	3 (16%)	22,30,33	1.86	5 (22%)
6	OMG	L5	1580	6	23,26,27	1.20	3 (13%)	33,38,41	1.93	10 (30%)

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
6	A2M	L5	3557	6	-	0/9/27/28	0/3/3/3
6	OMU	L5	2680	6	-	0/9/27/28	0/2/2/2
52	OMC	S2	174	52	-	0/9/27/28	0/2/2/2
52	PSU	S2	682	52	-	0/7/25/26	0/2/2/2
52	OMU	S2	121	52	-	0/9/27/28	0/2/2/2
6	PSU	L5	2351	6	-	0/7/25/26	0/2/2/2
6	OMC	L5	2194	6	-	1/9/27/28	0/2/2/2
52	OMU	S2	355	71,52	-	0/9/27/28	0/2/2/2
6	A2M	L5	398	6	-	0/9/27/28	0/3/3/3
6	OMC	L5	3433	6,90	-	4/9/27/28	0/2/2/2
52	OMG	S2	510	52,89	-	2/9/27/28	0/3/3/3
6	OMG	L5	1260	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	1720	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	823	52	-	2/7/25/26	0/2/2/2
6	PSU	L5	4246	6	-	1/7/25/26	0/2/2/2
6	PSU	L5	4740	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	3490	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	1638	6,90	-	0/7/25/26	0/2/2/2
52	PSU	S2	610	52	-	0/7/25/26	0/2/2/2
6	OMC	L5	2208	6,89	-	0/9/27/28	0/2/2/2
6	PSU	L5	3447	6	-	0/7/25/26	0/2/2/2
6	OMC	L5	3601	6	-	0/9/27/28	0/2/2/2
6	PSU	L5	4435	6	-	0/7/25/26	0/2/2/2
52	A2M	S2	669	52,89	-	2/9/27/28	0/3/3/3
6	PSU	L5	1718	6	-	0/7/25/26	0/2/2/2
47	MLZ	Lq	53	47	-	0/7/8/10	-
6	PSU	L5	3652	6,89	-	0/7/25/26	0/2/2/2
52	PSU	S2	1005	52	-	1/7/25/26	0/2/2/2
52	PSU	S2	1178	52	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PSU	L5	3371	6	-	0/7/25/26	0/2/2/2
52	OMG	S2	602	52	-	0/9/27/28	0/3/3/3
6	PSU	L5	3585	6,89	-	0/7/25/26	0/2/2/2
6	OMG	L5	4138	6	-	0/9/27/28	0/3/3/3
6	OMG	L5	2719	6	-	1/9/27/28	0/3/3/3
6	OMG	L5	4116	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	4217	6	-	0/7/25/26	0/2/2/2
52	OMG	S2	1491	52,89	-	1/9/27/28	0/3/3/3
6	PSU	L5	3554	6	-	0/7/25/26	0/2/2/2
6	A2M	L5	3562	6	-	0/9/27/28	0/3/3/3
6	A2M	L5	400	6	-	0/9/27/28	0/3/3/3
52	PSU	S2	1082	52	-	0/7/25/26	0/2/2/2
52	OMU	S2	1443	52,89	-	0/9/27/28	0/2/2/2
6	OMU	L5	4244	6	-	0/9/27/28	0/2/2/2
6	OMG	L5	4364	6	-	0/9/27/28	0/3/3/3
52	PSU	S2	407	52	-	0/7/25/26	0/2/2/2
6	OMG	L5	4383	6,90	-	0/9/27/28	0/3/3/3
52	OMG	S2	645	52	-	3/9/27/28	0/3/3/3
52	PSU	S2	1057	52	-	0/7/25/26	0/2/2/2
6	OMU	L5	2258	6	-	0/9/27/28	0/2/2/2
6	A2M	L5	1270	6	-	1/9/27/28	0/3/3/3
52	PSU	S2	119	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	1239	52	-	0/7/25/26	0/2/2/2
11	AYA	LF	2	11	-	2/4/6/8	-
6	PSU	L5	4169	6	-	0/7/25/26	0/2/2/2
52	A2M	S2	591	52	-	1/9/27/28	0/3/3/3
35	MLZ	Le	5	35	-	2/7/8/10	-
52	PSU	S2	864	52	-	0/7/25/26	0/2/2/2
52	OMU	S2	172	52	-	0/9/27/28	0/2/2/2
52	OMC	S2	518	52	-	2/9/27/28	0/2/2/2
52	MA6	S2	1852	52	-	2/11/29/30	0/3/3/3
46	M3L	Lp	98	46	-	0/9/10/12	-
6	A2M	L5	3517	6	-	3/9/27/28	0/3/3/3
6	OMU	L5	4052	6	-	0/9/27/28	0/2/2/2
52	OMU	S2	429	52	-	4/9/27/28	0/2/2/2
6	OMC	L5	4202	6	-	0/9/27/28	0/2/2/2
8	PSU	L8	55	8	-	0/7/25/26	0/2/2/2
52	MA6	S2	1851	52	-	2/11/29/30	0/3/3/3
6	PSU	L5	3466	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	1491	6	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
52	PSU	S2	816	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	3583	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	4039	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	4045	6	-	0/7/25/26	0/2/2/2
6	6MZ	L5	3966	6	-	0/9/27/28	0/3/3/3
6	OMC	L5	3619	6,22	-	2/9/27/28	0/2/2/2
6	PSU	L5	3500	6	-	0/7/25/26	0/2/2/2
52	OMU	S2	628	52	-	6/9/27/28	0/2/2/2
6	OMU	L5	4366	6	-	0/9/27/28	0/2/2/2
52	4AC	S2	1338	52	-	4/11/29/30	0/2/2/2
52	OMU	S2	1327	52,89	-	0/9/27/28	0/2/2/2
52	PSU	S2	218	52	-	0/7/25/26	0/2/2/2
6	OMG	L5	3524	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	4099	6	-	0/7/25/26	0/2/2/2
6	OMG	L5	3942	6	-	0/9/27/28	0/3/3/3
52	OMG	S2	1448	52	-	3/9/27/28	0/3/3/3
6	PSU	L5	3502	6	-	0/7/25/26	0/2/2/2
52	A2M	S2	469	52	-	0/9/27/28	0/3/3/3
6	A2M	L5	2244	6	-	1/9/27/28	0/3/3/3
52	PSU	S2	36	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	4382	6	-	2/7/25/26	0/2/2/2
6	A2M	L5	3450	6	-	0/9/27/28	0/3/3/3
6	OMC	L5	2667	6	-	0/9/27/28	0/2/2/2
6	OMG	L5	3676	6	-	1/9/27/28	0/3/3/3
6	PSU	L5	4322	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	4267	6,89	-	0/7/25/26	0/2/2/2
52	OMU	S2	1805	52	-	0/9/27/28	0/2/2/2
6	PSU	L5	3427	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	34	52	-	0/7/25/26	0/2/2/2
52	6MZ	S2	1833	52,89	-	1/9/27/28	0/3/3/3
6	PSU	L5	3496	6	-	0/7/25/26	0/2/2/2
6	OMG	L5	4369	6	-	2/9/27/28	0/3/3/3
6	A2M	L5	3599	6	-	0/9/27/28	0/3/3/3
6	OMU	L5	3973	6	-	0/9/27/28	0/2/2/2
52	PSU	S2	802	52	-	3/7/25/26	0/2/2/2
6	PSU	L5	4058	6	-	0/7/25/26	0/2/2/2
52	A2M	S2	159	52	-	0/9/27/28	0/3/3/3
52	PSU	S2	1245	52	-	0/7/25/26	0/2/2/2
81	AME	Sg	1	81	-	2/9/10/12	-
52	PSU	S2	687	52	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	OMC	L5	2265	6,89	-	2/9/27/28	0/2/2/2
6	PSU	L5	1537	6	-	0/7/25/26	0/2/2/2
6	OMC	L5	2647	6	-	0/9/27/28	0/2/2/2
6	PSU	L5	3462	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	3369	6,90,89	-	0/7/25/26	0/2/2/2
52	A2M	S2	577	52	-	1/9/27/28	0/3/3/3
83	HY3	Si	62	83	-	1/1/12/14	0/1/1/1
6	OMG	L5	3631	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	1632	6	-	1/7/25/26	0/2/2/2
52	PSU	S2	1047	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	1731	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	3576	6	-	1/7/25/26	0/2/2/2
6	A2M	L5	4336	6	-	2/9/27/28	0/3/3/3
52	A2M	S2	1384	52	-	0/9/27/28	0/3/3/3
6	PSU	L5	3494	6	-	2/7/25/26	0/2/2/2
10	HIC	LE	245	10	-	2/5/6/8	0/1/1/1
52	PSU	S2	1368	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	1046	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	4374	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	1644	52,89	-	0/7/25/26	0/2/2/2
52	PSU	S2	93	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	109	52	-	0/7/25/26	0/2/2/2
52	A2M	S2	485	52	-	1/9/27/28	0/3/3/3
6	A2M	L5	3492	6,89	-	3/9/27/28	0/3/3/3
52	PSU	S2	1693	52	-	0/7/25/26	0/2/2/2
6	OMU	L5	3657	6	-	0/9/27/28	0/2/2/2
6	5MC	L5	4193	6	-	3/7/25/26	0/2/2/2
52	OMC	S2	1392	52	-	2/9/27/28	0/2/2/2
52	PSU	S2	652	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	1348	52	-	0/7/25/26	0/2/2/2
52	OMU	S2	116	52	-	0/9/27/28	0/2/2/2
52	PSU	S2	1233	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	573	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	4166	6	-	2/7/25/26	0/2/2/2
6	PSU	L5	4177	6	-	0/7/25/26	0/2/2/2
52	OMC	S2	1704	52,89	-	0/9/27/28	0/2/2/2
6	OMG	L5	3476	6	-	0/9/27/28	0/3/3/3
6	OMG	L5	4240	6	-	1/9/27/28	0/3/3/3
52	A2M	S2	27	52	-	0/9/27/28	0/3/3/3
6	UR3	L5	4276	6	-	0/7/25/26	0/2/2/2

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	OMC	L5	3540	6	-	0/9/27/28	0/2/2/2
52	OMG	S2	684	52	-	2/9/27/28	0/3/3/3
52	OMG	S2	868	52	-	0/9/27/28	0/3/3/3
6	OMG	L5	1477	6	-	3/9/27/28	0/3/3/3
6	PSU	L5	4203	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	105	52	-	0/7/25/26	0/2/2/2
6	OMC	L5	1820	6,89	-	0/9/27/28	0/2/2/2
6	A2M	L5	3456	6	-	0/9/27/28	0/3/3/3
52	A2M	S2	1679	52	-	0/9/27/28	0/3/3/3
6	PSU	L5	1799	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	1801	6	-	0/7/25/26	0/2/2/2
6	A2M	L5	2658	6,89	-	1/9/27/28	0/3/3/3
6	PSU	L5	4149	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	1446	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	1721	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	4188	6	-	0/7/25/26	0/2/2/2
6	PSU	L5	4107	6	-	0/7/25/26	0/2/2/2
6	OMC	L5	2704	6	-	0/9/27/28	0/2/2/2
6	PSU	L5	4711	6	-	0/7/25/26	0/2/2/2
6	1MA	L5	1266	6,89	-	2/7/25/26	0/3/3/3
6	A2M	L5	1810	6,89	-	0/9/27/28	0/3/3/3
6	OMG	L5	3359	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	4749	6	-	0/7/25/26	0/2/2/2
6	A2M	L5	2630	6	-	3/9/27/28	0/3/3/3
52	A2M	S2	513	52	-	0/9/27/28	0/3/3/3
52	PSU	S2	650	52	-	0/7/25/26	0/2/2/2
52	PSU	S2	1626	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	4419	6	-	0/7/25/26	0/2/2/2
52	A2M	S2	1032	52	-	0/9/27/28	0/3/3/3
8	PSU	L8	69	8	-	0/7/25/26	0/2/2/2
6	A2M	L5	4317	6	-	0/9/27/28	0/3/3/3
52	OMC	S2	463	52	-	0/9/27/28	0/2/2/2
6	OMC	L5	3573	6	-	0/9/27/28	0/2/2/2
52	OMU	S2	1289	52,55	-	0/9/27/28	0/2/2/2
52	A2M	S2	99	52,89	-	1/9/27/28	0/3/3/3
6	5MC	L5	3514	6,89	-	0/7/25/26	0/2/2/2
6	OMG	L5	4245	6	-	0/9/27/28	0/3/3/3
52	PSU	S2	815	52	-	0/7/25/26	0/2/2/2
6	PSU	L5	4325	6	-	0/7/25/26	0/2/2/2
52	7MG	S2	1640	52	-	0/7/37/38	0/3/3/3
6	OMG	L5	2267	6	-	0/9/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
52	PSU	S2	967	52	-	0/7/25/26	0/2/2/2
52	A2M	S2	166	52	-	0/9/27/28	0/3/3/3
6	PSU	L5	1683	6,90	-	0/7/25/26	0/2/2/2
6	OMG	L5	3974	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	2475	6	-	0/7/25/26	0/2/2/2
6	OMC	L5	4282	6	-	0/9/27/28	0/2/2/2
6	A2M	L5	4269	6,89	-	0/9/27/28	0/3/3/3
52	4AC	S2	1843	52,86	-	2/11/29/30	0/2/2/2
8	OMG	L8	75	8	-	0/9/27/28	0/3/3/3
52	PSU	S2	867	52	-	0/7/25/26	0/2/2/2
6	A2M	L5	1479	6	-	2/9/27/28	0/3/3/3
6	OMG	L5	2207	6	-	0/9/27/28	0/3/3/3
6	PSU	L5	4278	6	-	0/7/25/26	0/2/2/2
52	PSU	S2	210	52	-	0/7/25/26	0/2/2/2
49	SAC	Ls	2	49	-	0/7/8/10	-
6	PSU	L5	4298	6	-	0/7/25/26	0/2/2/2
78	SAC	Sd	2	78	-	0/7/8/10	-
52	PSU	S2	1175	52	-	0/7/25/26	0/2/2/2
52	OMG	S2	1329	52	-	0/9/27/28	0/3/3/3
79	NMM	Se	67	79	-	3/9/11/13	-
6	A2M	L5	2206	6,89	-	0/9/27/28	0/3/3/3
6	OMC	L5	1284	6	-	0/9/27/28	0/2/2/2
6	PSU	L5	4042	6	-	0/7/25/26	0/2/2/2
52	OMG	S2	437	52	-	0/9/27/28	0/3/3/3
60	SAC	SL	2	60	-	1/7/8/10	-
6	A2M	L5	1489	6,89	-	2/9/27/28	0/3/3/3
6	PSU	L5	3616	6	-	0/7/25/26	0/2/2/2
6	OMG	L5	1580	6	-	1/9/27/28	0/3/3/3

The worst 5 of 716 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L5	1266	1MA	C2-N3	4.79	1.39	1.30
52	S2	1851	MA6	C5-C4	4.52	1.47	1.39
52	S2	1852	MA6	C5-C4	4.44	1.47	1.39
52	S2	1833	6MZ	C5-C4	4.31	1.47	1.39
79	Se	67	NMM	CZ-NH2	4.23	1.44	1.34

The worst 5 of 1372 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	S2	1005	PSU	C4-N3-C2	-8.37	114.28	126.34
79	Se	67	NMM	NE-CZ-NH2	-7.51	112.59	119.48
6	L5	1266	1MA	N1-C2-N3	-6.95	117.73	126.00
52	S2	485	A2M	N3-C2-N1	-6.79	117.97	128.60
6	L5	2630	A2M	N3-C2-N1	-6.79	117.99	128.60

There are no chirality outliers.

5 of 111 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	LF	2	AYA	OT-CT-N-CA
11	LF	2	AYA	CM-CT-N-CA
35	Le	5	MLZ	N-CA-CB-CG
35	Le	5	MLZ	C-CA-CB-CG
52	S2	429	OMU	C2'-C1'-N1-C2

There are no ring outliers.

67 monomers are involved in 82 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	L5	3557	A2M	1	0
6	L5	2680	OMU	1	0
52	S2	682	PSU	1	0
6	L5	2194	OMC	2	0
6	L5	398	A2M	1	0
52	S2	610	PSU	1	0
6	L5	4138	OMG	1	0
6	L5	2719	OMG	2	0
6	L5	400	A2M	1	0
6	L5	4364	OMG	1	0
6	L5	4383	OMG	1	0
6	L5	1270	A2M	1	0
52	S2	1239	PSU	1	0
35	Le	5	MLZ	1	0
52	S2	172	OMU	1	0
52	S2	1852	MA6	1	0
6	L5	3517	A2M	1	0
6	L5	4202	OMC	1	0
52	S2	1851	MA6	1	0
6	L5	3466	PSU	1	0
6	L5	4366	OMU	2	0
52	S2	1338	4AC	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
52	S2	218	PSU	1	0
6	L5	3942	OMG	1	0
52	S2	1448	OMG	1	0
6	L5	4382	PSU	2	0
6	L5	3676	OMG	1	0
52	S2	1833	6MZ	1	0
6	L5	4369	OMG	2	0
6	L5	3599	A2M	2	0
6	L5	3973	OMU	2	0
81	Sg	1	AME	2	0
6	L5	2265	OMC	1	0
83	Si	62	HY3	1	0
6	L5	3631	OMG	1	0
6	L5	1632	PSU	1	0
52	S2	1047	PSU	2	0
52	S2	1046	PSU	1	0
6	L5	4374	PSU	1	0
52	S2	109	PSU	1	0
6	L5	3657	OMU	1	0
6	L5	4193	5MC	2	0
52	S2	116	OMU	2	0
52	S2	1233	PSU	2	0
52	S2	1704	OMC	1	0
6	L5	3476	OMG	1	0
6	L5	3540	OMC	1	0
6	L5	1477	OMG	2	0
6	L5	4203	PSU	1	0
52	S2	1679	A2M	1	0
6	L5	3359	OMG	1	0
52	S2	513	A2M	1	0
6	L5	4419	PSU	1	0
52	S2	1032	A2M	1	0
6	L5	3573	OMC	1	0
6	L5	3514	5MC	1	0
52	S2	1640	7MG	1	0
6	L5	4282	OMC	1	0
52	S2	1843	4AC	1	0
8	L8	75	OMG	1	0
6	L5	2207	OMG	2	0
6	L5	4278	PSU	1	0
6	L5	1284	OMC	1	0
52	S2	437	OMG	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
60	SL	2	SAC	1	0
6	L5	1489	A2M	1	0
6	L5	1580	OMG	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 256 ligands modelled in this entry, 253 are monoatomic - leaving 3 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	$\# Z > 2$
88	SER	F	101	2	4,5,6	0.57	0	0,5,7	-	-
87	GTP	E	1001	-	30,34,34	0.84	1 (3%)	46,54,54	1.68	10 (21%)
87	GTP	L7	203	7	30,34,34	1.11	4 (13%)	46,54,54	1.81	9 (19%)

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
88	SER	F	101	2	-	0/2/4/6	-
87	GTP	E	1001	-	-	6/22/38/38	0/3/3/3
87	GTP	L7	203	7	-	0/22/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
87	L7	203	GTP	C5-N7	-2.60	1.34	1.39
87	L7	203	GTP	C5-C4	2.54	1.45	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
87	L7	203	GTP	C6-N1	-2.42	1.34	1.38
87	L7	203	GTP	C4-N9	-2.21	1.32	1.38
87	E	1001	GTP	C2-N3	2.15	1.38	1.33

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
87	L7	203	GTP	C5-C4-N3	-5.40	119.70	128.46
87	E	1001	GTP	C5-C4-N3	-4.96	120.41	128.46
87	L7	203	GTP	N9-C4-N3	4.95	135.87	125.94
87	L7	203	GTP	C2-N3-C4	4.69	120.66	112.30
87	E	1001	GTP	C2-N3-C4	4.41	120.15	112.30

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

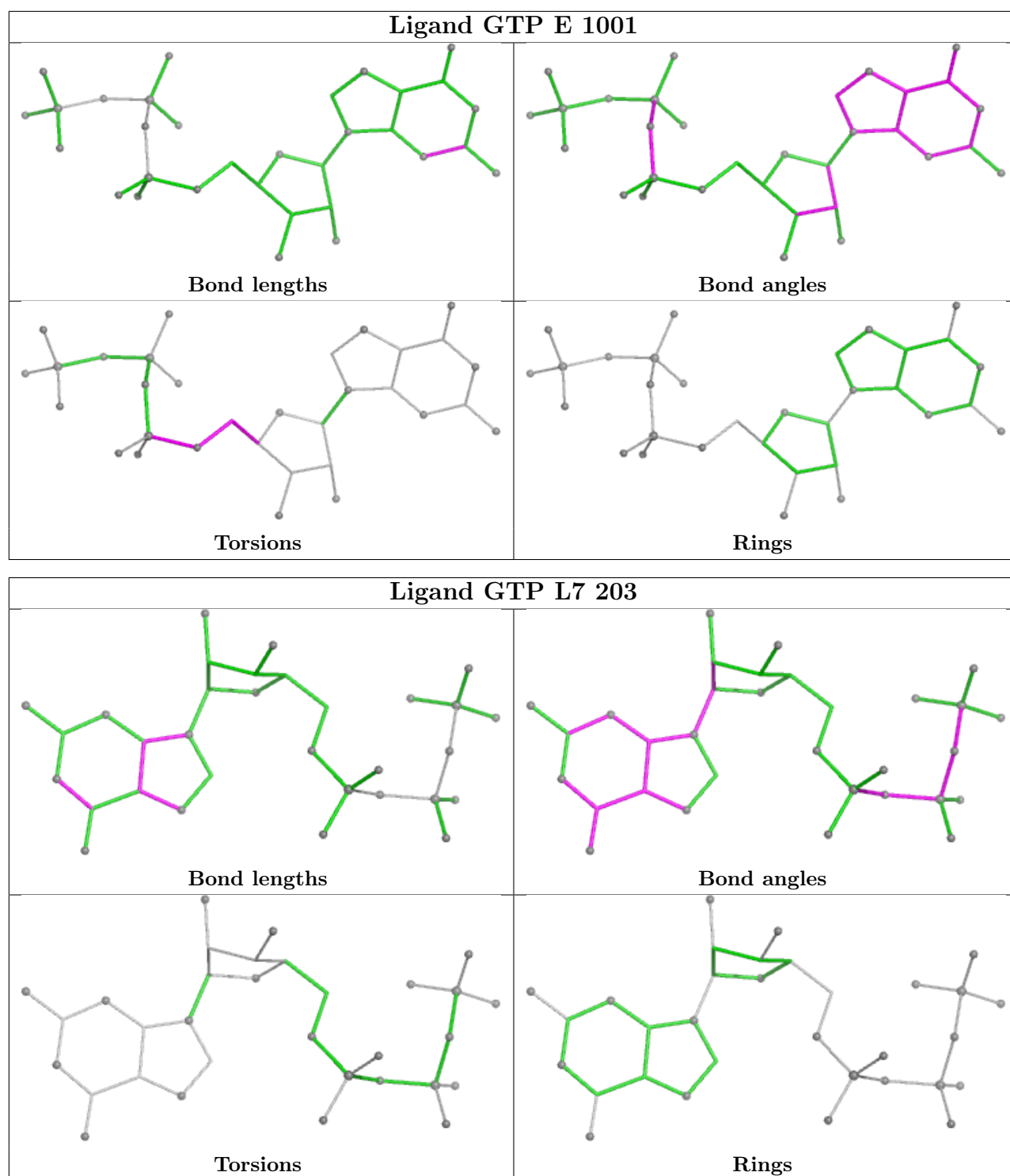
Mol	Chain	Res	Type	Atoms
87	E	1001	GTP	C5'-O5'-PA-O3A
87	E	1001	GTP	O4'-C4'-C5'-O5'
87	E	1001	GTP	C3'-C4'-C5'-O5'
87	E	1001	GTP	C5'-O5'-PA-O1A
87	E	1001	GTP	C5'-O5'-PA-O2A

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
87	E	1001	GTP	1	0
87	L7	203	GTP	3	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 ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
83	Si	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	Si	61:GLN	C	62:HY3	N	3.10

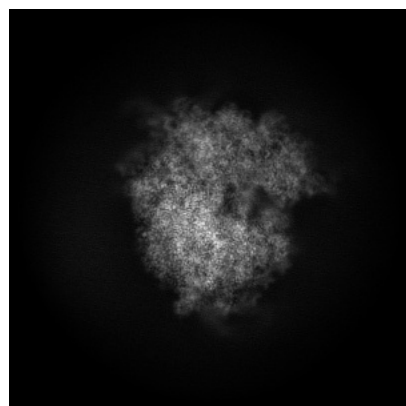
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14751. These allow visual inspection of the internal detail of the map and identification of artifacts.

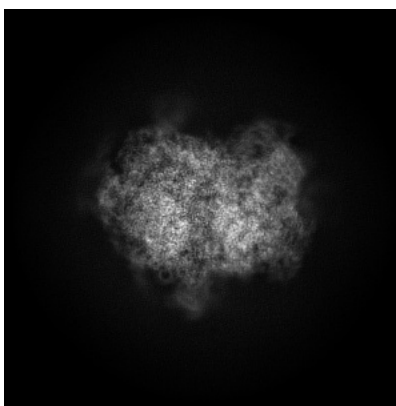
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

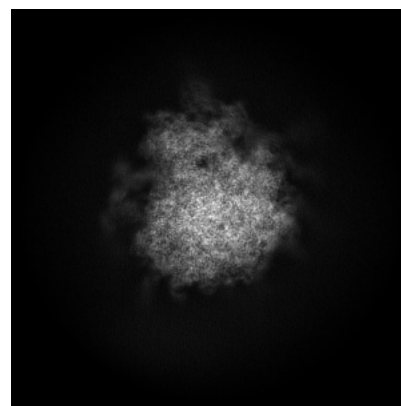
6.1.1 Primary map



X

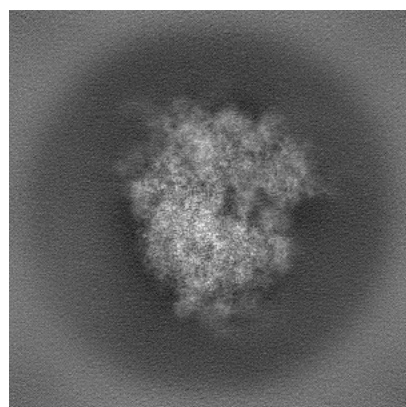


Y

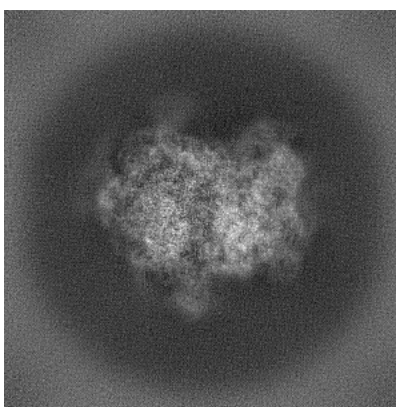


Z

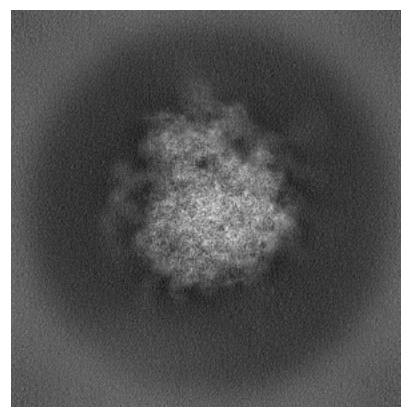
6.1.2 Raw map



X



Y

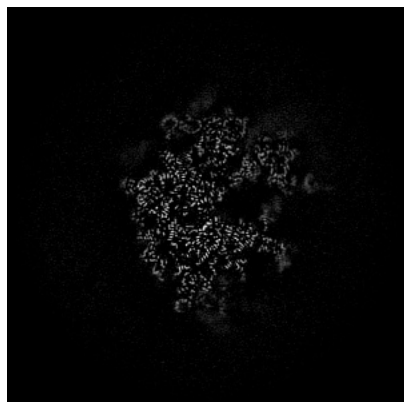


Z

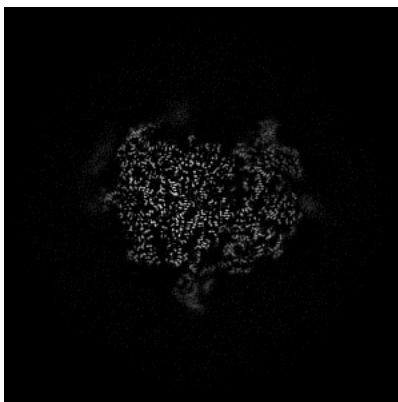
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

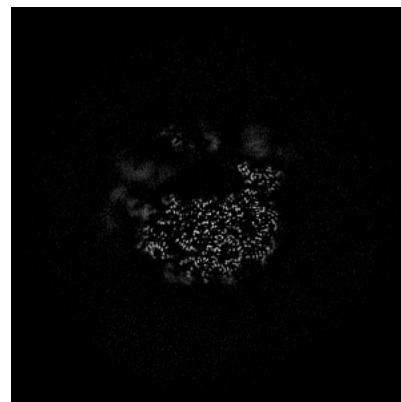
6.2.1 Primary map



X Index: 256

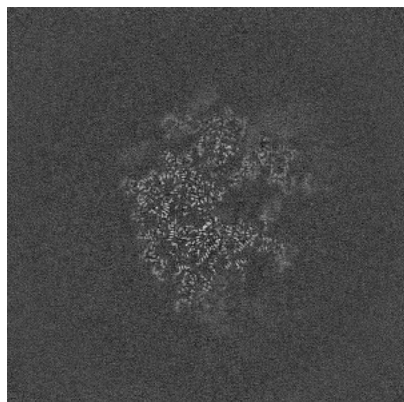


Y Index: 256

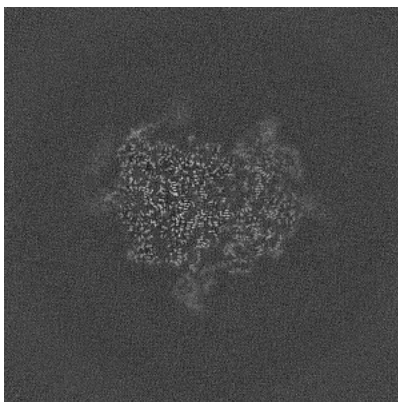


Z Index: 256

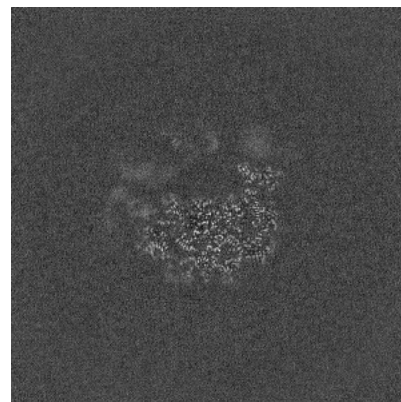
6.2.2 Raw map



X Index: 256



Y Index: 256

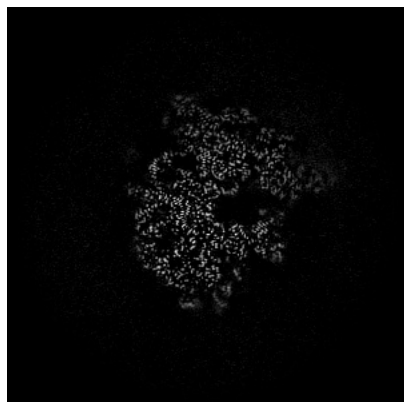


Z Index: 256

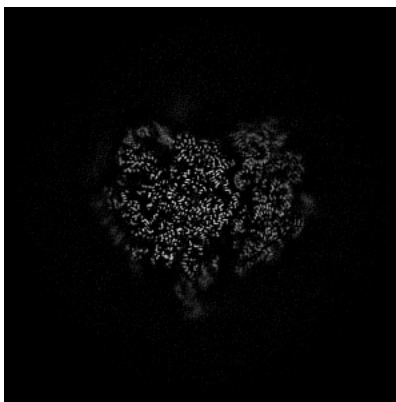
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

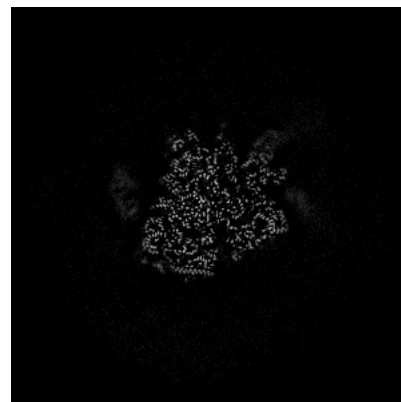
6.3.1 Primary map



X Index: 240

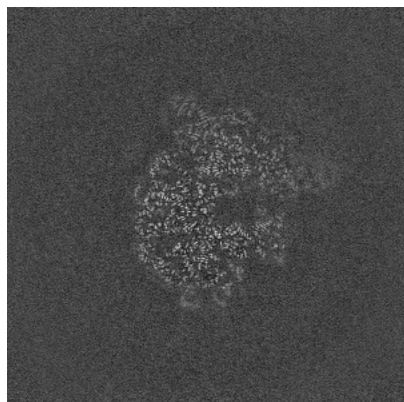


Y Index: 246

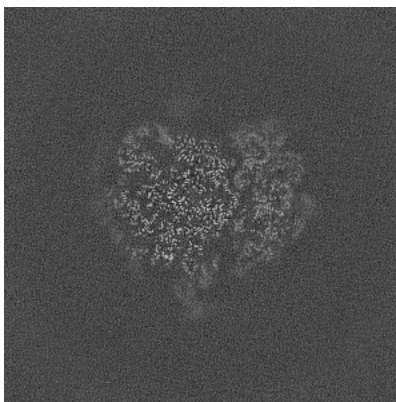


Z Index: 223

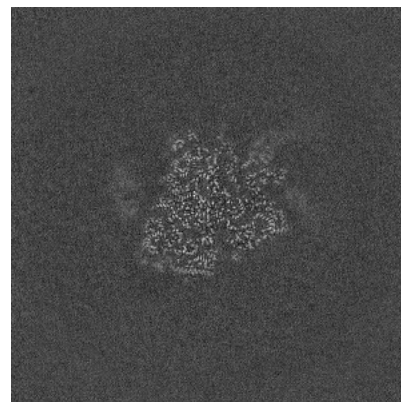
6.3.2 Raw map



X Index: 238



Y Index: 246

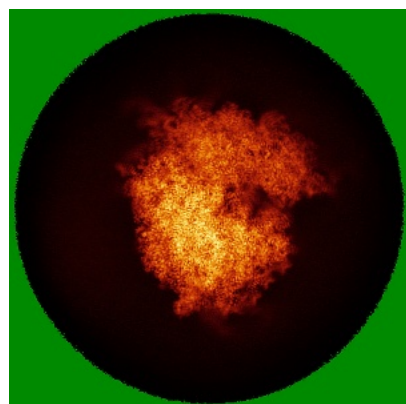


Z Index: 223

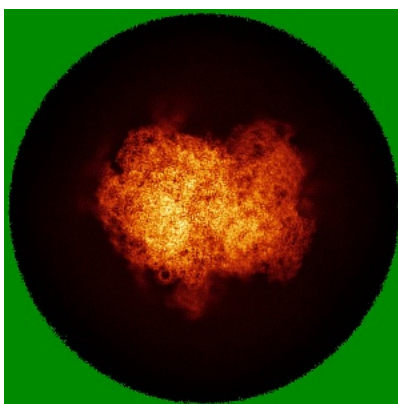
The images above show the largest variance slices of the map in three orthogonal directions.

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

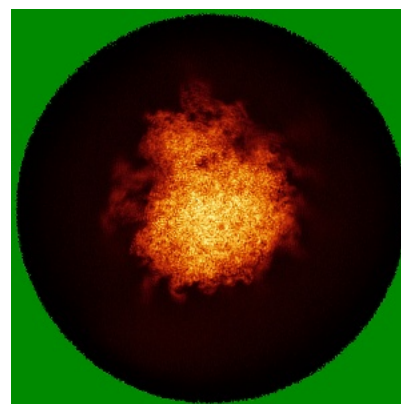
6.4.1 Primary map



X

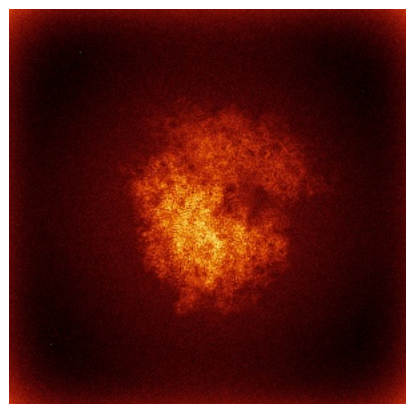


Y

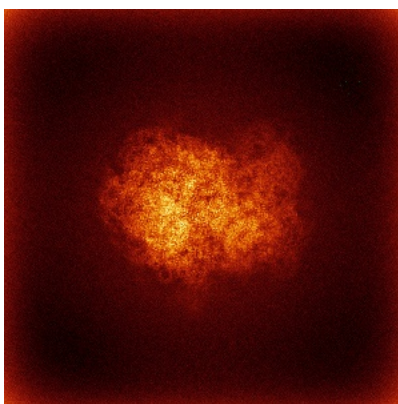


Z

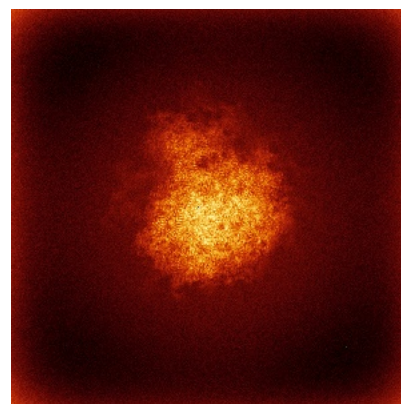
6.4.2 Raw map



X



Y

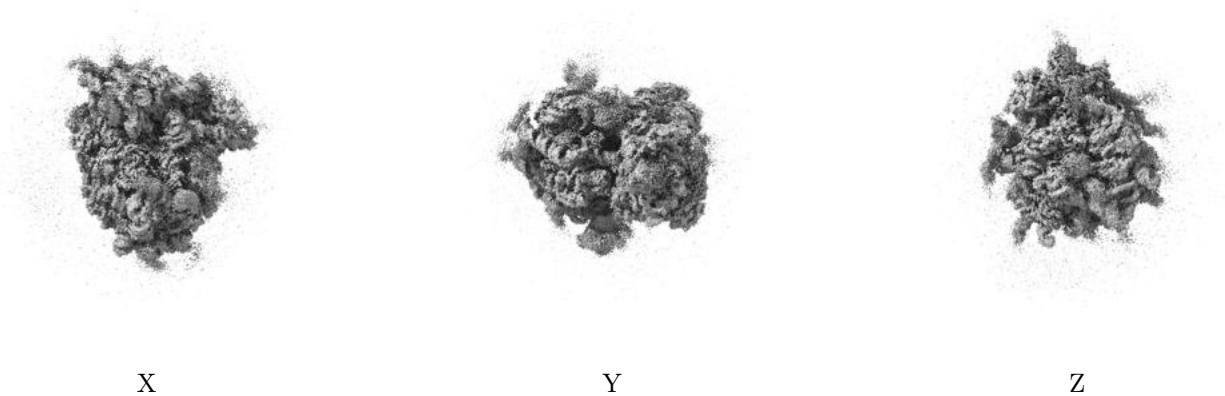


Z

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.

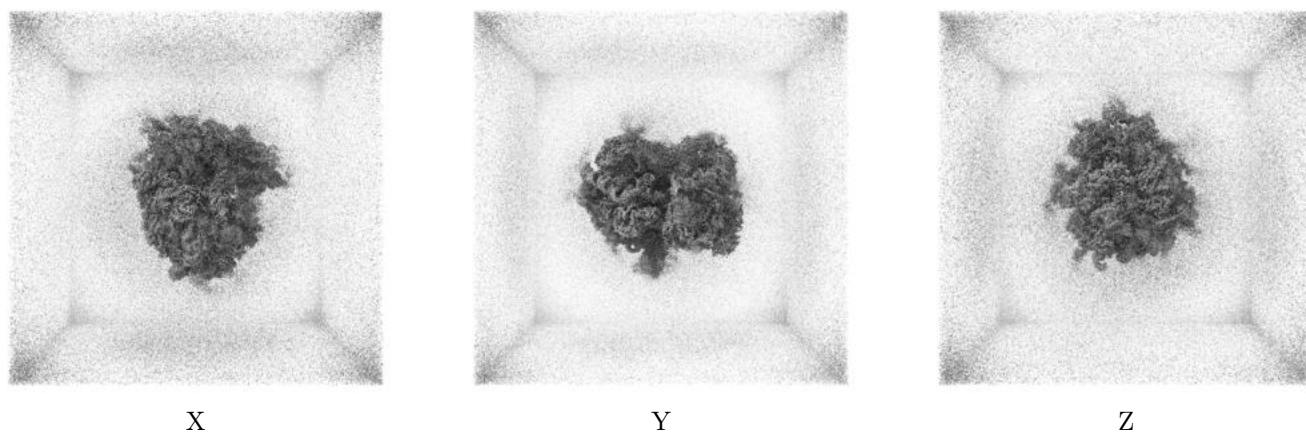
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.3. 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



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.

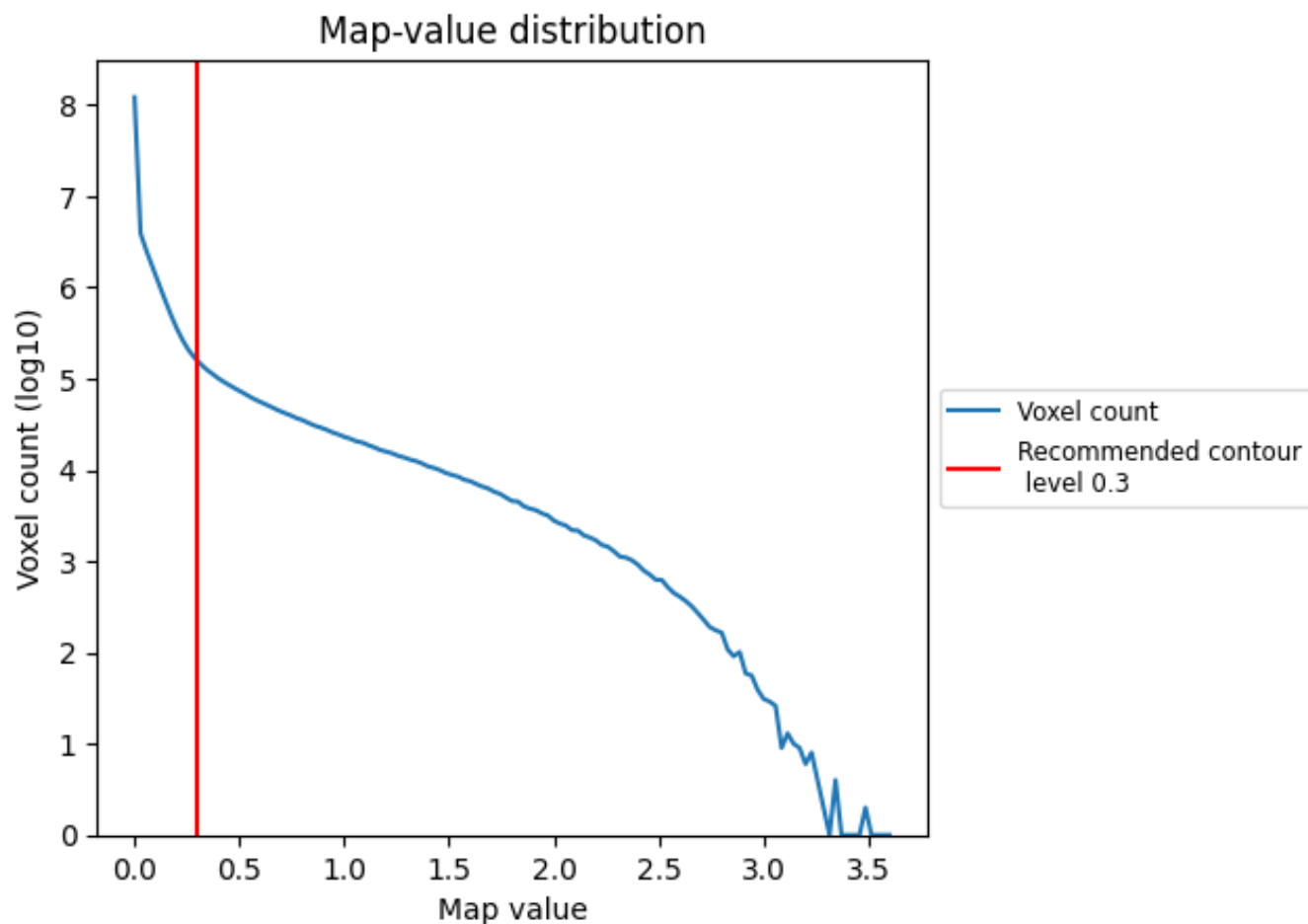
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

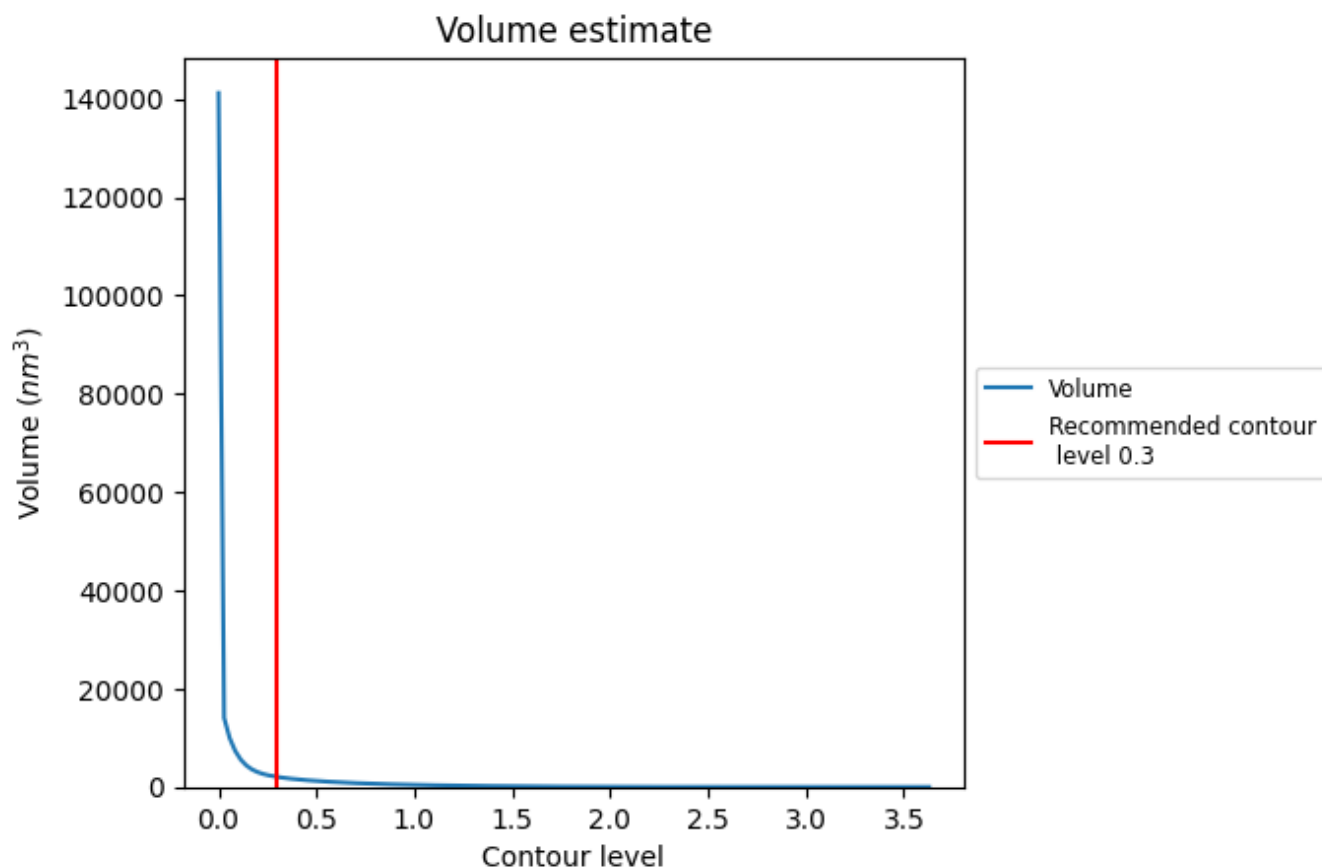
This section contains the results of statistical analysis of the map.

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.

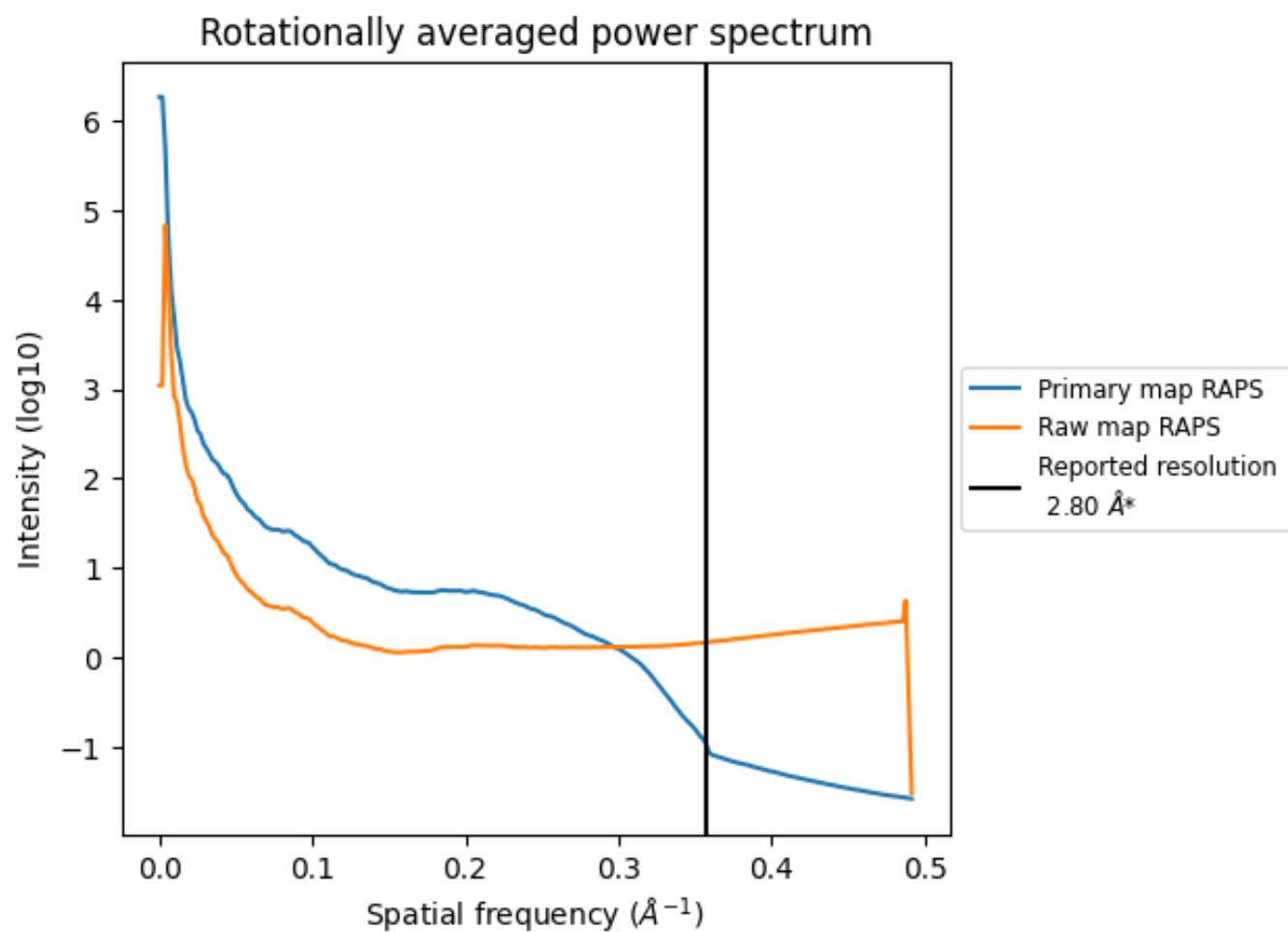
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2046 nm^3 ; this corresponds to an approximate mass of 1848 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 ⓘ

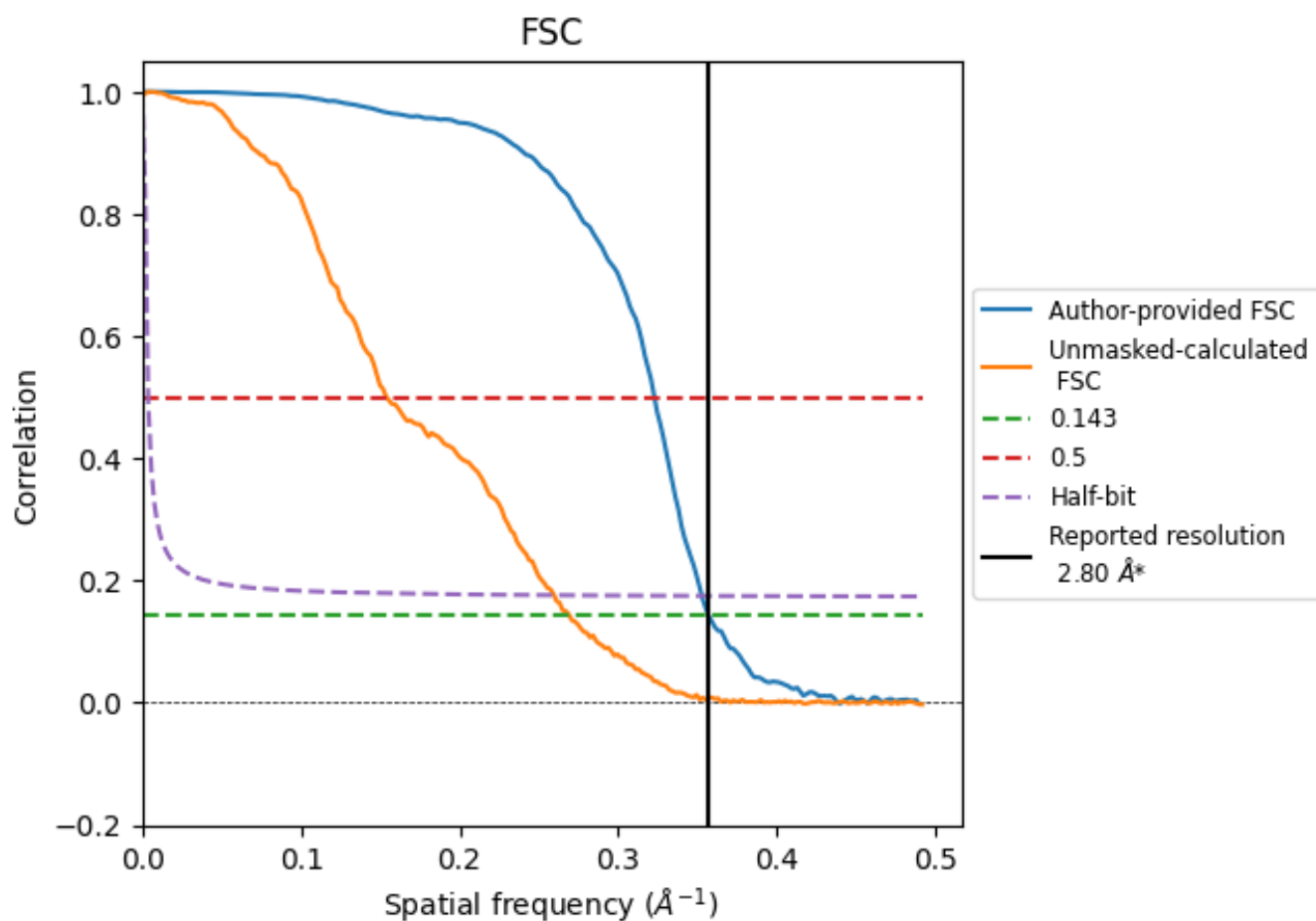


*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

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 0.357 Å⁻¹

8.2 Resolution estimates [i](#)

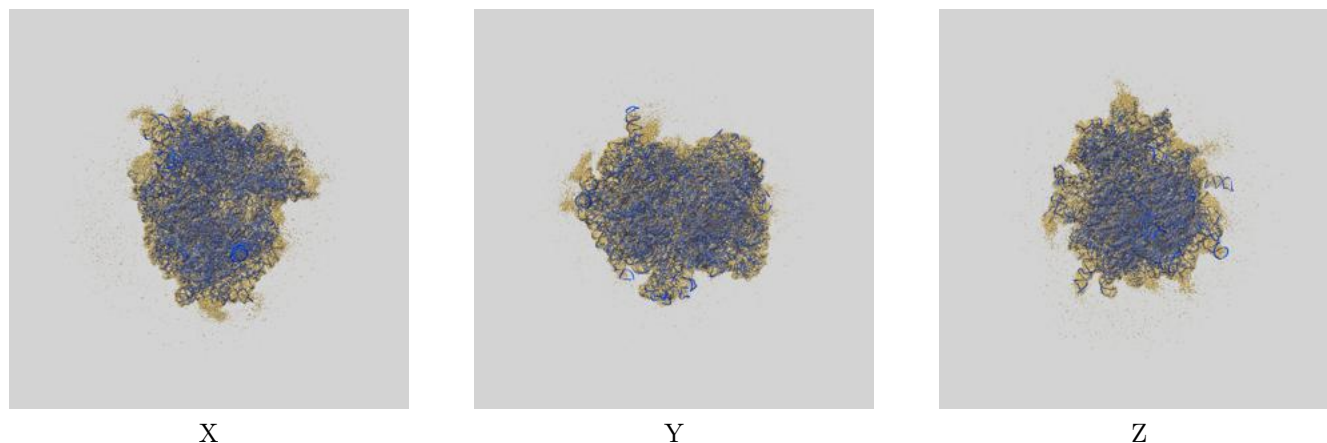
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	2.80	3.10	2.83
Unmasked-calculated*	3.72	6.49	3.85

*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 3.72 differs from the reported value 2.8 by more than 10 %

9 Map-model fit [i](#)

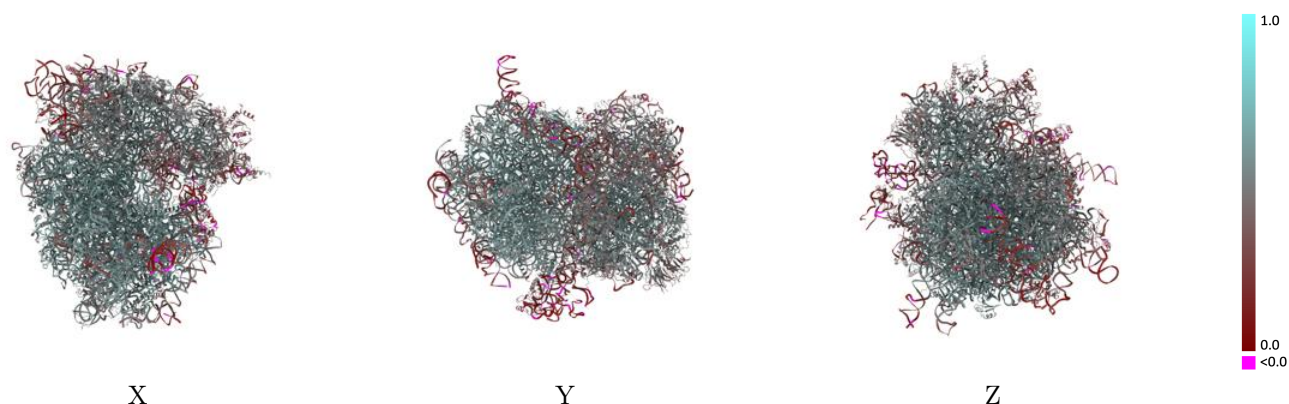
This section contains information regarding the fit between EMDB map EMD-14751 and PDB model 7ZJW. Per-residue inclusion information can be found in section [3](#) on page [25](#).

9.1 Map-model overlay [i](#)



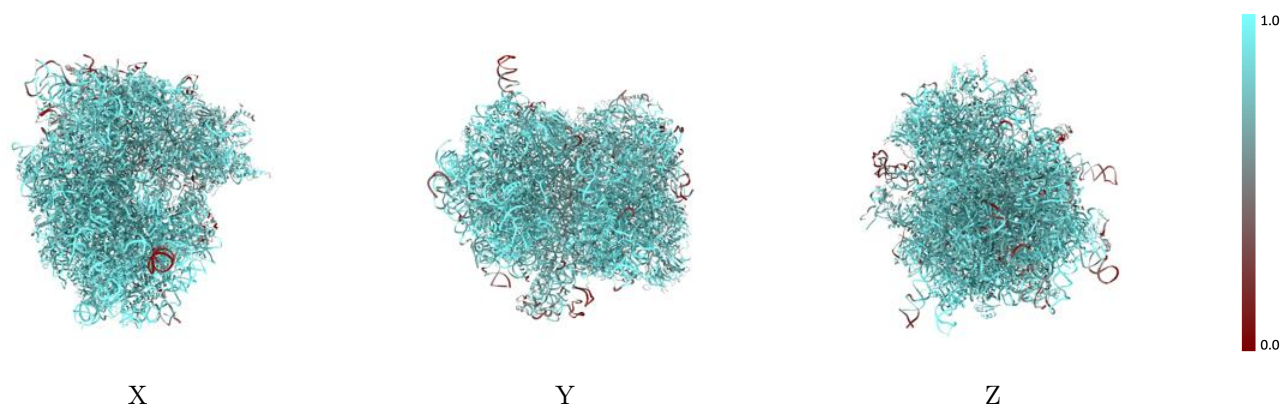
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



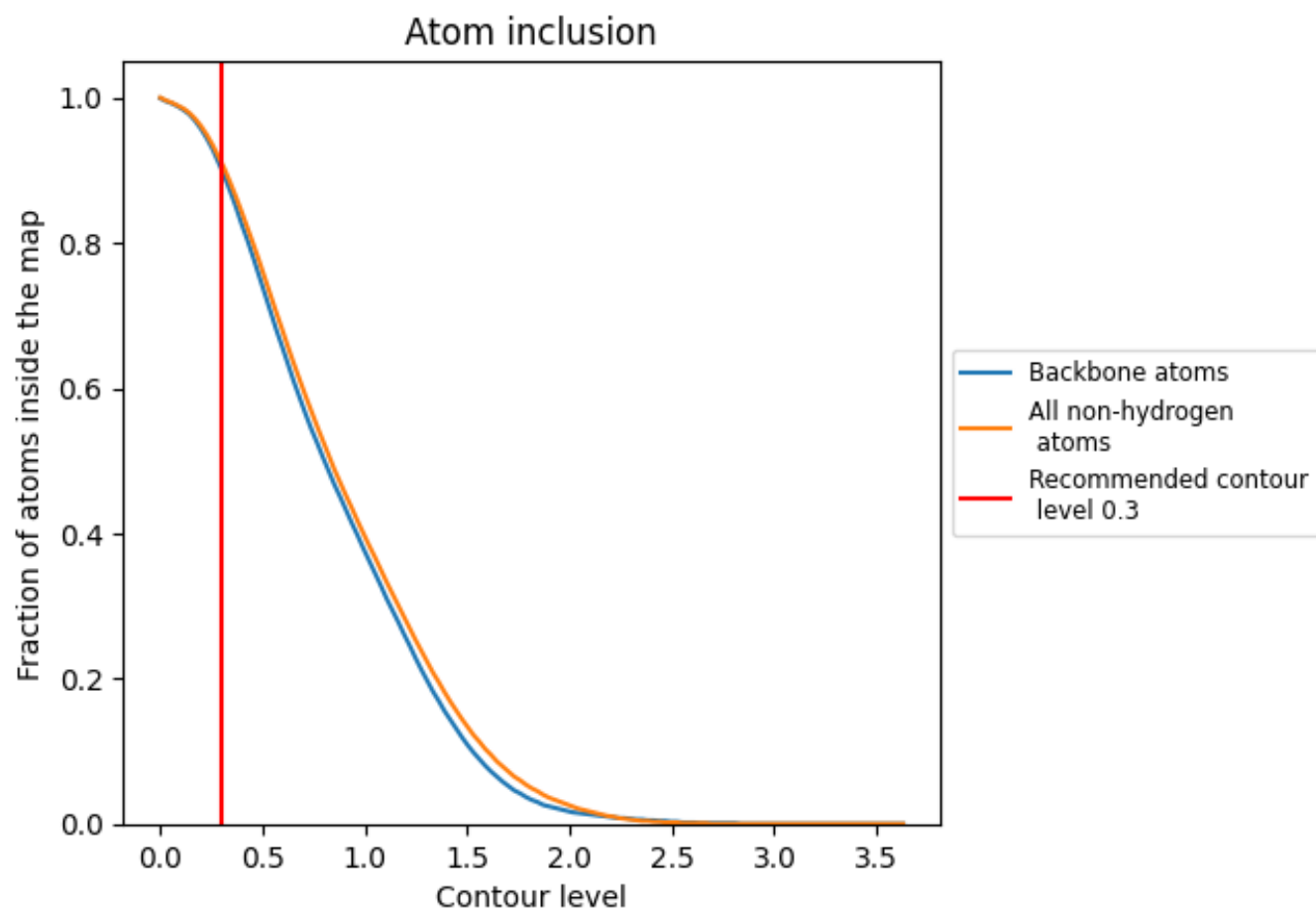
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

9.4 Atom inclusion ⓘ



At the recommended contour level, 90% of all backbone atoms, 91% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9100	 0.5140
B	 0.8170	 0.4050
E	 0.7650	 0.3700
F	 0.7450	 0.3280
I	 0.6340	 0.2040
L5	 0.9540	 0.5340
L7	 0.9970	 0.5910
L8	 0.9790	 0.5740
LD	 0.9440	 0.5950
LE	 0.9460	 0.5830
LF	 0.9480	 0.5850
LG	 0.9410	 0.5540
LH	 0.9140	 0.5460
LI	 0.9430	 0.5850
LJ	 0.8990	 0.5380
LK	 0.9230	 0.5660
LL	 0.9030	 0.5590
LM	 0.9110	 0.5300
LO	 0.9160	 0.5610
LP	 0.9340	 0.5650
LQ	 0.9700	 0.6080
LR	 0.9490	 0.5870
LS	 0.9220	 0.5770
LT	 0.9480	 0.5860
LU	 0.9150	 0.5480
LV	 0.9610	 0.5940
LW	 0.9200	 0.5650
LX	 0.9210	 0.5110
LY	 0.9000	 0.5750
LZ	 0.8160	 0.4910
La	 0.9320	 0.5610
Lb	 0.9310	 0.5730
Lc	 0.9430	 0.5590
Ld	 0.9630	 0.5940
Le	 0.8330	 0.4890



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
Lf	 0.8520	 0.5250
Lg	 0.9360	 0.5700
Lh	 0.9430	 0.5910
Li	 0.9520	 0.5940
Lj	 0.9280	 0.5670
Lk	 0.9250	 0.5620
Ll	 0.9250	 0.5570
Lm	 0.9670	 0.6070
Ln	 0.8760	 0.5360
Lo	 0.9370	 0.5810
Lp	 0.9310	 0.5710
Lq	 0.9190	 0.5790
Lr	 0.9270	 0.5760
Ls	 0.9540	 0.5850
Lt	 0.6190	 0.2220
Lx	 0.6990	 0.2950
S	 0.8890	 0.2700
S2	 0.9330	 0.4940
SB	 0.8230	 0.5070
SC	 0.8070	 0.4900
SD	 0.6820	 0.3590
SE	 0.7160	 0.4180
SF	 0.8980	 0.5410
SG	 0.8230	 0.4210
SH	 0.9050	 0.5370
SL	 0.8620	 0.5140
SM	 0.8600	 0.5110
SN	 0.8820	 0.5300
SO	 0.8300	 0.4860
SP	 0.8270	 0.4790
SQ	 0.8430	 0.4820
SR	 0.8170	 0.4060
SS	 0.7670	 0.4550
ST	 0.8380	 0.4760
SU	 0.8260	 0.4680
SV	 0.8490	 0.4690
SW	 0.7760	 0.4980
SX	 0.6610	 0.3040
SY	 0.8800	 0.5410
SZ	 0.8760	 0.5310
Sa	 0.7780	 0.4480
Sb	 0.8620	 0.5000

Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
Sc	 0.8050	 0.4680
Sd	 0.8620	 0.4840
Se	 0.8730	 0.4940
Sf	 0.8160	 0.4630
Sg	 0.8670	 0.5210
Sh	 0.9120	 0.5560
Si	 0.8980	 0.5470
Sj	 0.8500	 0.4590
Sk	 0.8080	 0.4450
Sl	 0.9270	 0.5970