



Full wwPDB EM Validation Report ⓘ

Jul 2, 2024 – 03:54 PM JST

PDB ID : 8XSY
EMDB ID : EMD-38630
Title : Cryo-EM structure of the human 80S ribosome with Tigecycline, e-tRNA and CCDC124 (40S head Swivelled)
Authors : Li, X.; Wang, M.; Cheng, J.
Deposited on : 2024-01-10
Resolution : 3.00 Å(reported)
Based on initial model : 6Z6M

This is a Full wwPDB EM Validation 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:

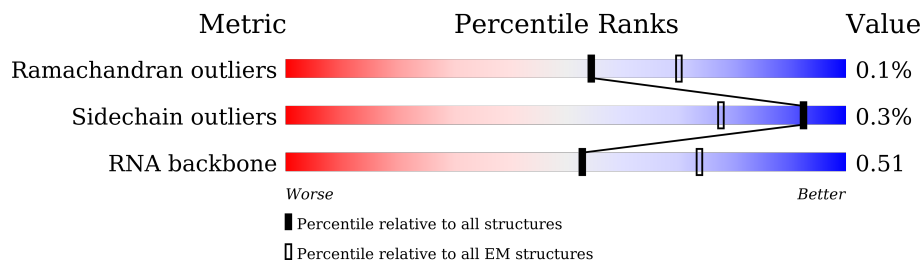
EMDB validation analysis : 0.0.1.dev92
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

1 Overall quality at a glance

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

The reported resolution of this entry is 3.00 Å.

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



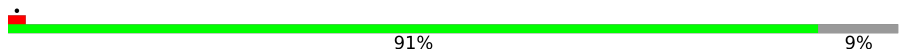
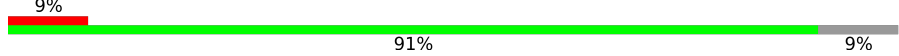
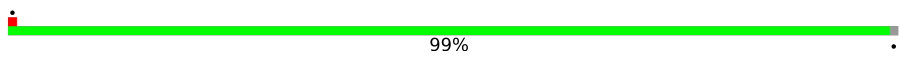
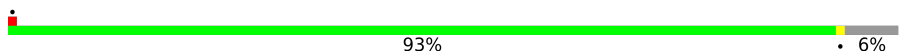
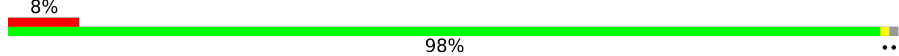
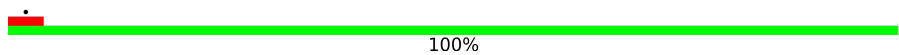

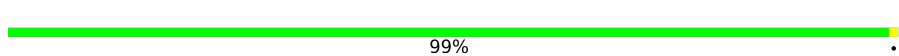
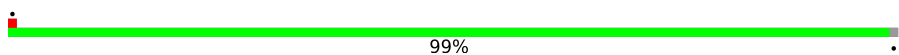

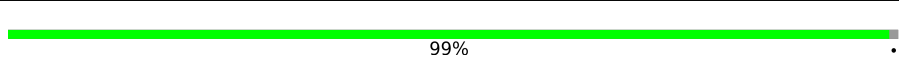
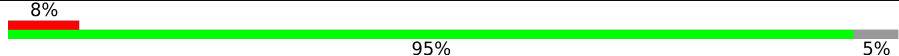
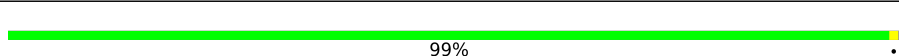
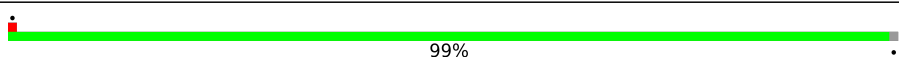
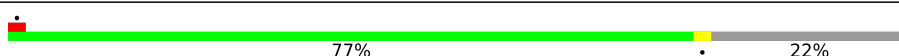
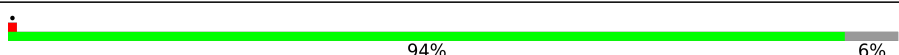
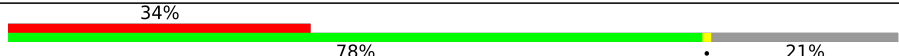
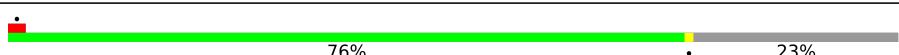
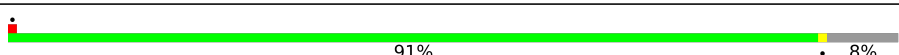
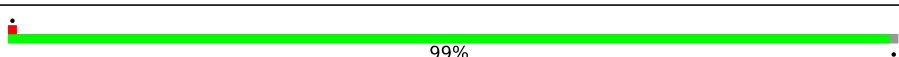
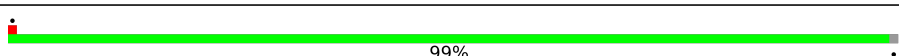



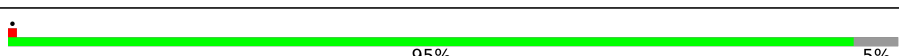
Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	L5	5070	
2	L7	121	
3	L8	157	
4	LA	257	
5	LB	403	
6	LC	427	
7	LD	297	
8	LE	288	

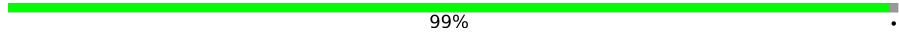
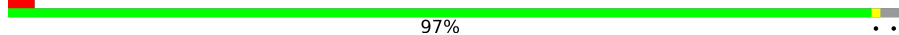
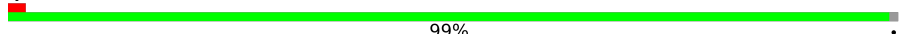
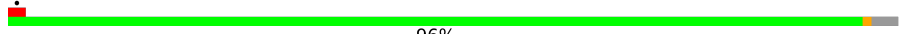


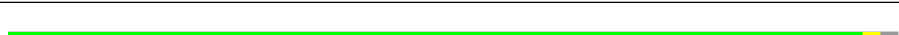
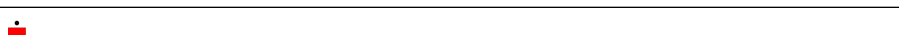
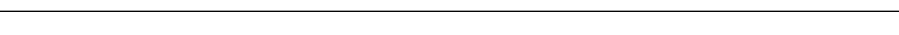
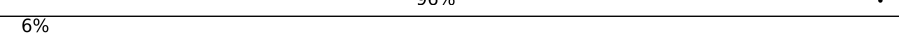
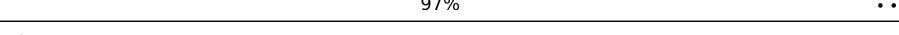
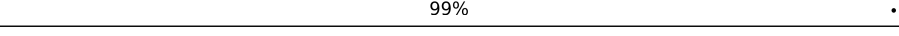











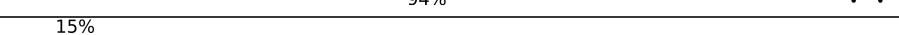
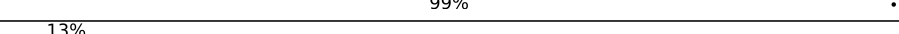
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Mol	Chain	Length	Quality of chain
9	LF	248	 91% 9%
10	LG	266	 91% 9%
11	LH	192	 99%
12	LI	214	 93% 6%
13	LJ	178	 98%
14	LL	211	 100%
15	LM	215	 65% 35%
16	LN	204	 99%
17	LO	203	 99%
18	LP	184	 82% 17%
19	LQ	188	 99%
20	LR	196	 95% 5%
21	LS	176	 99%
22	LT	160	 99%
23	LU	128	 77% 22%
24	LV	140	 94% 6%
25	LW	157	 78% 21%
26	LX	156	 76% 23%
27	LY	145	 91% 8%
28	LZ	136	 99%
29	La	148	 99%
30	Lb	159	 68% 31%
31	Lc	115	 85% 15%
32	Ld	125	 86% 14%
33	Le	135	 95% 5%

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Mol	Chain	Length	Quality of chain
34	Lf	110	 99%
35	Lg	117	 97%
36	Lh	123	 99%
37	Li	105	 96%
38	Lj	97	 89% 11%
39	Lk	70	 99%
40	Ll	51	 96%
41	Lm	128	 41% 59%
42	Ln	25	 96%
43	Lo	106	 6% 97%
44	Lp	92	 99%
45	Lr	137	 91% 9%
46	Ls	317	 58% 62% 38%
47	Lt	165	 84% 84% 15%
48	Lz	217	 94% 94% 6%
49	S2	1869	 7% 67% 26% 7%
50	SA	295	 6% 74% 25%
51	SB	264	 8% 80% 19%
52	SD	243	 18% 93% 7%
53	SE	263	 10% 99%
54	SF	204	 10% 89% 10%
55	SH	194	 18% 94%
56	SI	208	 15% 99%
57	SK	165	 13% 57% 42%
58	SL	158	 9% 91% 9%

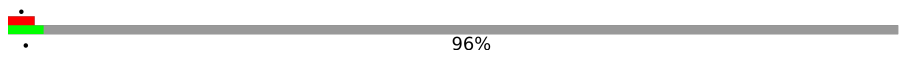


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Mol	Chain	Length	Quality of chain
59	SP	145	22% 88% 11%
60	SQ	146	12% 98%
61	SR	135	10% 98%
62	SS	152	21% 93% 5%
63	ST	145	13% 98%
64	SU	119	19% 87% 13%
65	SV	83	6% 100%
66	SX	143	98%
67	Sa	115	87% 11%
68	Sc	69	22% 93% 7%
69	Sd	56	11% 96%
70	Sg	317	17% 98%
71	SC	293	75% 24%
72	SG	249	20% 93% 5%
73	SJ	194	11% 95%
74	SM	132	70% 92% 8%
75	SN	151	99%
76	SO	151	8% 89% 11%
77	SW	130	98%
78	SY	133	17% 94%
79	SZ	125	18% 57% 40%
80	Sb	84	12% 99%
81	Se	59	25% 97%
82	Sf	156	30% 40% 57%
83	CA	394	90% 89% 10%

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Mol	Chain	Length	Quality of chain
84	CB	408	 96%
85	CC	75	 68% 69% 23% 8%
86	CE	223	 33% 33% 67%

2 Entry composition [i](#)

There are 89 unique types of molecules in this entry. The entry contains 223822 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 28S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	L5	3771	80096	35636	14582	26108	3770	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L5	2113	C	G	conflict	GB 86475748

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	L7	120	2561	1141	456	844	120	0	0

- Molecule 3 is a RNA chain called 5.8S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
3	L8	156	3314	1480	585	1094	155	0	0

- Molecule 4 is a protein called 60S ribosomal protein L8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	LA	248	1898	1189	389	314	6	0	0

- Molecule 5 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	LB	402	3238	2060	608	556	14	0	0

- Molecule 6 is a protein called 60S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	LC	368	2927	1840	583	489	15	0	0

- Molecule 7 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	LD	293	2382	1507	434	427	14	0	0

- Molecule 8 is a protein called 60S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	LE	221	1774	1142	336	292	4	0	0

- Molecule 9 is a protein called 60S ribosomal protein L7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	LF	225	1870	1202	358	301	9	0	0

- Molecule 10 is a protein called 60S ribosomal protein L7a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	LG	241	1927	1228	371	324	4	0	0

- Molecule 11 is a protein called 60S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	LH	190	1518	956	284	272	6	0	0

- Molecule 12 is a protein called Large ribosomal subunit protein uL16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	LI	202	1639	1041	316	269	13	0	0

- Molecule 13 is a protein called 60S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	LJ	176	Total	C	N	O	S	0	0
			1410	888	263	253	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
14	LL	210	Total	C	N	O	S	0	0
			1701	1064	352	281	4		

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
17	LO	201	Total	C	N	O	S	0	0
			1650	1063	321	261	5		

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
20	LR	187	1566	971	336	250	9	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
21	LS	175	1453	925	283	235	10	0	0

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
23	LU	100	816	524	142	148	2	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
24	LV	131	979	618	184	172	5	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
25	LW	124	1015	634	207	170	4	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
26	LX	120	985	630	185	169	1	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
27	LY	133	1106	694	224	185	3	0	0

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
30	Lb	109	882	549	192	137	4	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
31	Lc	98	764	485	135	138	6	0	0

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
33	Le	128	1053	667	216	165	5	0	0

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
37	Li	102	832	521	177	129	5	0	0

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
41	Lm	52	Total	C	N	O	S	0	0
			430	267	90	67	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
42	Ln	24	Total	C	N	O	S	0	0
			230	139	62	26	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
43	Lo	105	Total	C	N	O	S	0	0
			862	542	175	139	6		

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

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

- Molecule 45 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	Lr	125	Total	C	N	O	S	0	0
			1002	622	207	168	5		

- Molecule 46 is a protein called Large ribosomal subunit protein uL10.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	Ls	196	Total	C	N	O	S	0	0
			1496	952	259	276	9		

- Molecule 47 is a protein called 60S ribosomal protein L12.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	Lt	141	Total	C	N	O	S	0	0
			1046	652	191	199	4		

- Molecule 48 is a protein called 60S ribosomal protein L10a.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
48	Lz	205	1018	607	205	206	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
49	S2	1740	36896	16458	6597	12102	1739	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S2	582	C	U	conflict	GB 36162
S2	583	C	A	conflict	GB 36162
S2	584	G	A	conflict	GB 36162
S2	798	A	G	conflict	GB 36162
S2	1095	U	C	conflict	GB 36162

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
50	SA	221	1741	1106	305	322	8	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
51	SB	214	1738	1103	310	311	14	0	0

- Molecule 52 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
52	SD	227	1765	1125	317	315	8	0	0

- Molecule 53 is a protein called 40S ribosomal protein S4, X isoform.

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

- Molecule 54 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
54	SF	184	1461	914	276	264	7	0	0

- Molecule 55 is a protein called 40S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
55	SH	186	1497	956	274	266	1	0	0

- Molecule 56 is a protein called 40S ribosomal protein S8.

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

- Molecule 57 is a protein called 40S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
57	SK	95	799	524	139	130	6	0	0

- Molecule 58 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
58	SL	144	1182	752	224	200	6	0	0

- Molecule 59 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
59	SP	129	1061	672	202	180	7	0	0

- Molecule 60 is a protein called 40S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
60	SQ	144	1142	726	216	197	3	0	0

- Molecule 61 is a protein called 40S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
61	SR	135	1090	685	202	198	5	0	0

- Molecule 62 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
62	SS	144	1190	746	241	202	1	0	0

- Molecule 63 is a protein called 40S ribosomal protein S19.

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

- Molecule 64 is a protein called 40S ribosomal protein S20.

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

- Molecule 65 is a protein called 40S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
65	SV	83	636	393	117	121	5	0	0

- Molecule 66 is a protein called 40S ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
66	SX	141	1098	693	219	183	3	0	0

- Molecule 67 is a protein called 40S ribosomal protein S26.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
67	Sa	102	821	512	171	133	5	0	0

- Molecule 68 is a protein called 40S ribosomal protein S28.

Mol	Chain	Residues	Atoms					AltConf	Trace
68	Sc	64	Total	C	N	O	S	0	0
			506	308	102	94	2		

- Molecule 69 is a protein called 40S ribosomal protein S29.

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

- Molecule 70 is a protein called Receptor of activated protein C kinase 1.

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

- Molecule 71 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
71	SC	222	Total	C	N	O	S	0	0
			1725	1115	298	302	10		

- Molecule 72 is a protein called 40S ribosomal protein S6.

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

- Molecule 73 is a protein called 40S ribosomal protein S9.

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

- Molecule 74 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms				AltConf	Trace
74	SM	122	Total	C	N	O	0	0
			604	359	122	123		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
SM	52	GLN	LEU	conflict	UNP P25398
SM	69	LEU	CYS	conflict	UNP P25398
SM	99	ASN	LYS	conflict	UNP P25398

- Molecule 75 is a protein called 40S ribosomal protein S13.

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

- Molecule 76 is a protein called 40S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
76	SO	135	1010	618	198	188	6	0	0

- Molecule 77 is a protein called 40S ribosomal protein S15a.

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

- Molecule 78 is a protein called 40S ribosomal protein S24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
78	SY	126	1027	648	201	173	5	0	0

- Molecule 79 is a protein called 40S ribosomal protein S25.

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

- Molecule 80 is a protein called 40S ribosomal protein S27.

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

- Molecule 81 is a protein called 40S ribosomal protein S30.

Mol	Chain	Residues	Atoms					AltConf	Trace
81	Se	58	Total	C	N	O	S	0	0
			459	284	100	74	1		

- Molecule 82 is a protein called Ubiquitin-40S ribosomal protein S27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
82	Sf	67	Total	C	N	O	S	0	0
			548	346	102	93	7		

- Molecule 83 is a protein called Proliferation-associated protein 2G4.

Mol	Chain	Residues	Atoms					AltConf	Trace
83	CA	354	Total	C	N	O	S	4	0
			2764	1744	475	528	17		

- Molecule 84 is a protein called SERPINE1 mRNA-binding protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
84	CB	16	Total	C	N	O	S	0	0
			127	80	22	24	1		

- Molecule 85 is a RNA chain called E-tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
85	CC	69	Total	C	N	O	P	0	0
			1482	661	277	475	69		

- Molecule 86 is a protein called Coiled-coil domain-containing protein 124.

Mol	Chain	Residues	Atoms					AltConf	Trace
86	CE	73	Total	C	N	O	S	0	0
			613	369	122	121	1		

- Molecule 87 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

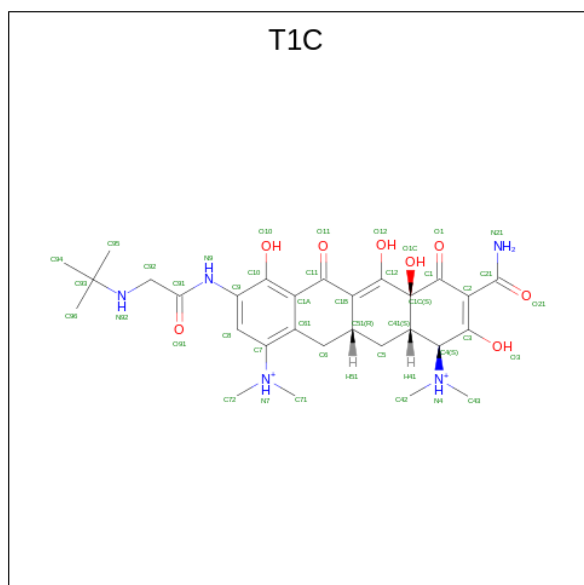
Mol	Chain	Residues	Atoms		AltConf
87	L5	212	Total	Mg	0
			212	212	
87	L7	3	Total	Mg	0
			3	3	
87	L8	4	Total	Mg	0
			4	4	

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Mol	Chain	Residues	Atoms		AltConf
87	LA	1	Total	Mg	0
			1	1	
87	LP	1	Total	Mg	0
			1	1	
87	LV	1	Total	Mg	0
			1	1	
87	Le	1	Total	Mg	0
			1	1	
87	Lg	1	Total	Mg	0
			1	1	
87	Lj	1	Total	Mg	0
			1	1	
87	S2	28	Total	Mg	0
			28	28	
87	SQ	1	Total	Mg	0
			1	1	
87	SG	1	Total	Mg	0
			1	1	

- Molecule 88 is TIGECYCLINE (three-letter code: T1C) (formula: C₂₉H₄₁N₅O₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf	
88	L5	1	Total	C	N	O	0
			42	29	5	8	
88	L5	1	Total	C	N	O	0
			42	29	5	8	

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
88	L5	1	42	29	5	8	0
88	L5	1	42	29	5	8	0
88	L5	1	42	29	5	8	0
88	L5	1	42	29	5	8	0
88	L5	1	42	29	5	8	0
88	CC	1	42	29	5	8	0

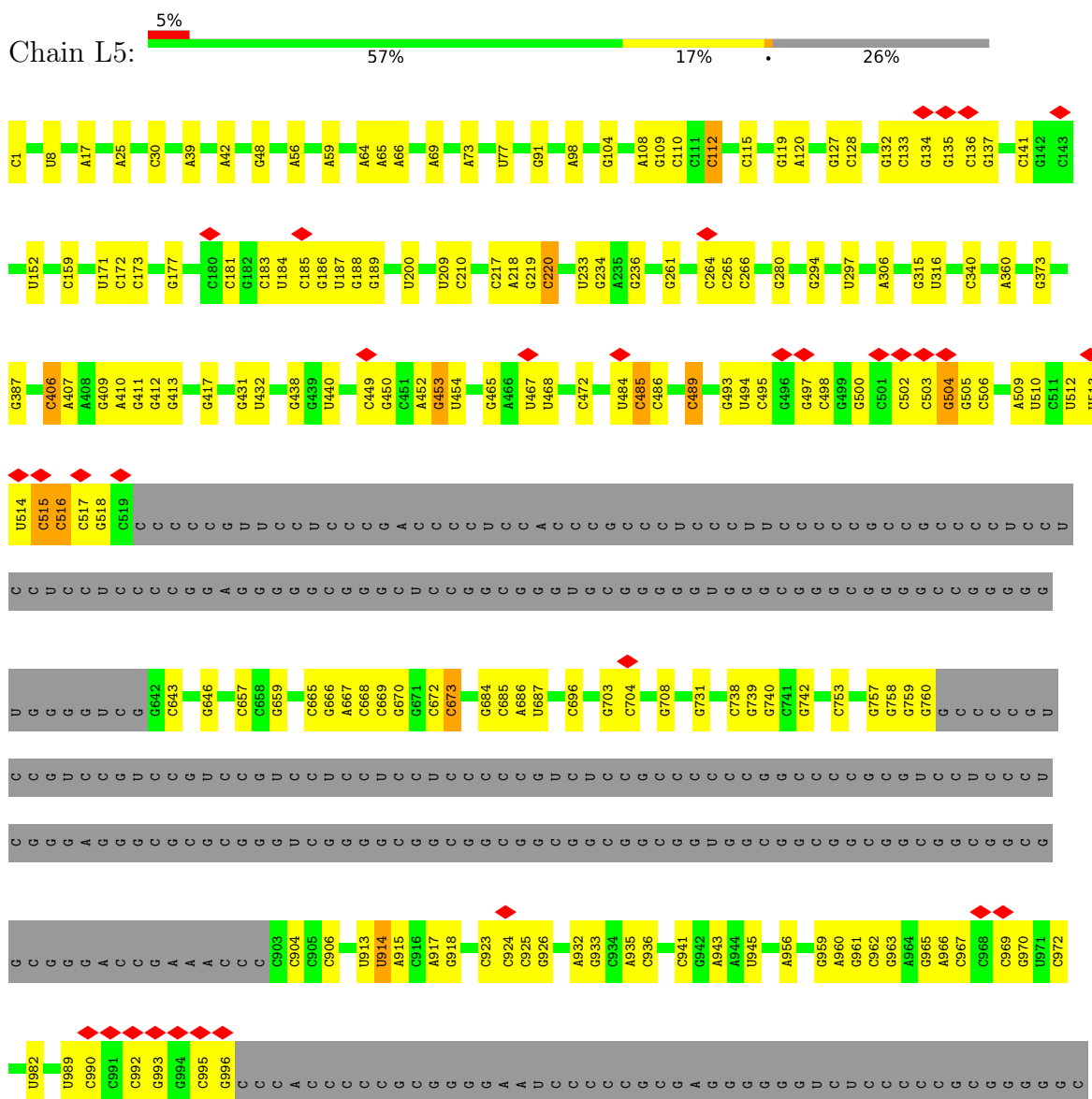
- Molecule 89 is ZINC ION (three-letter code: ZN) (formula: Zn).

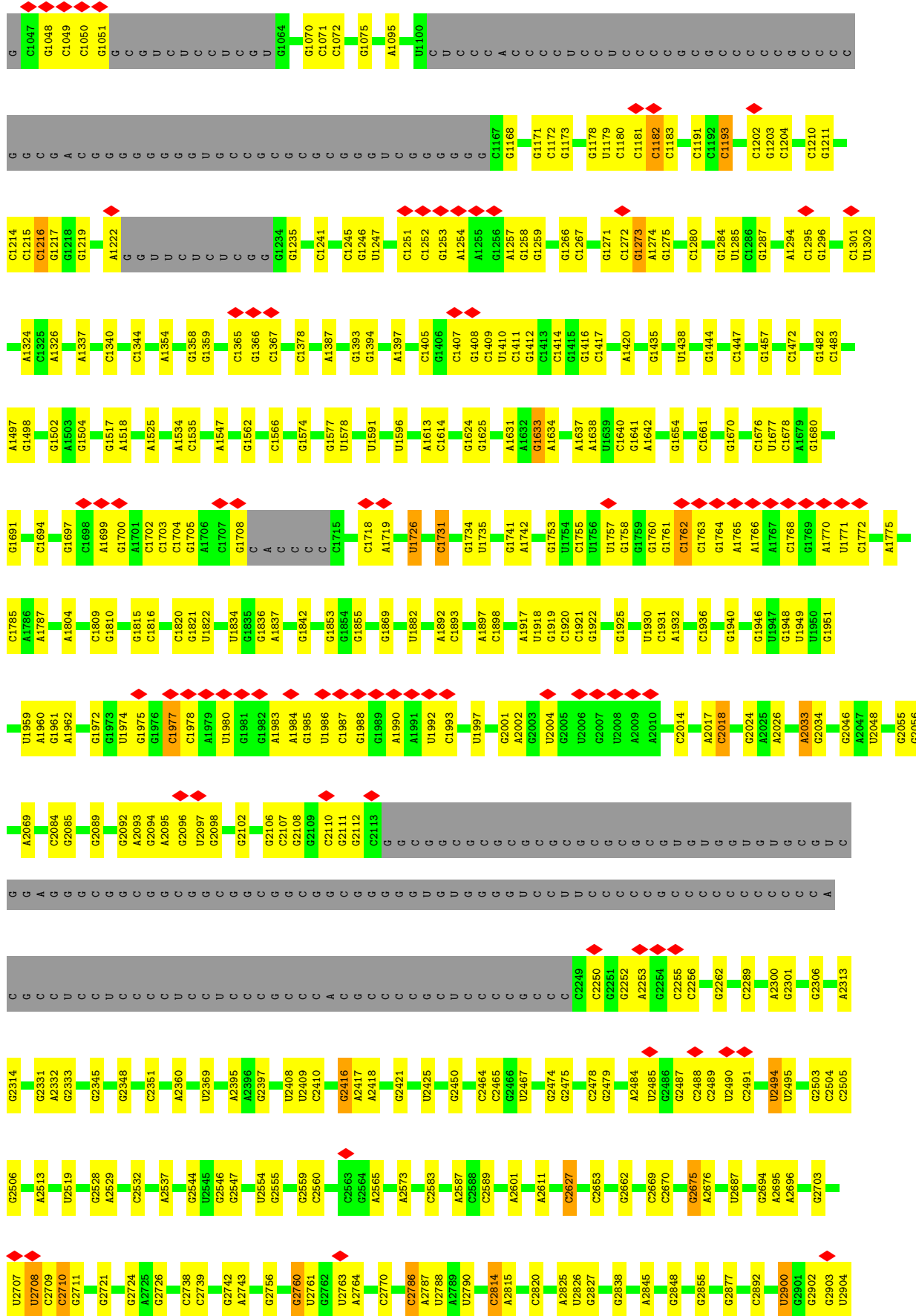
Mol	Chain	Residues	Atoms		AltConf
			Total	Zn	
89	Lg	1	1	1	0
89	Lj	1	1	1	0
89	Lm	1	1	1	0
89	Lo	1	1	1	0
89	Lp	1	1	1	0
89	SD	1	1	1	0
89	Sa	1	1	1	0
89	SM	1	1	1	0

3 Residue-property plots [i](#)

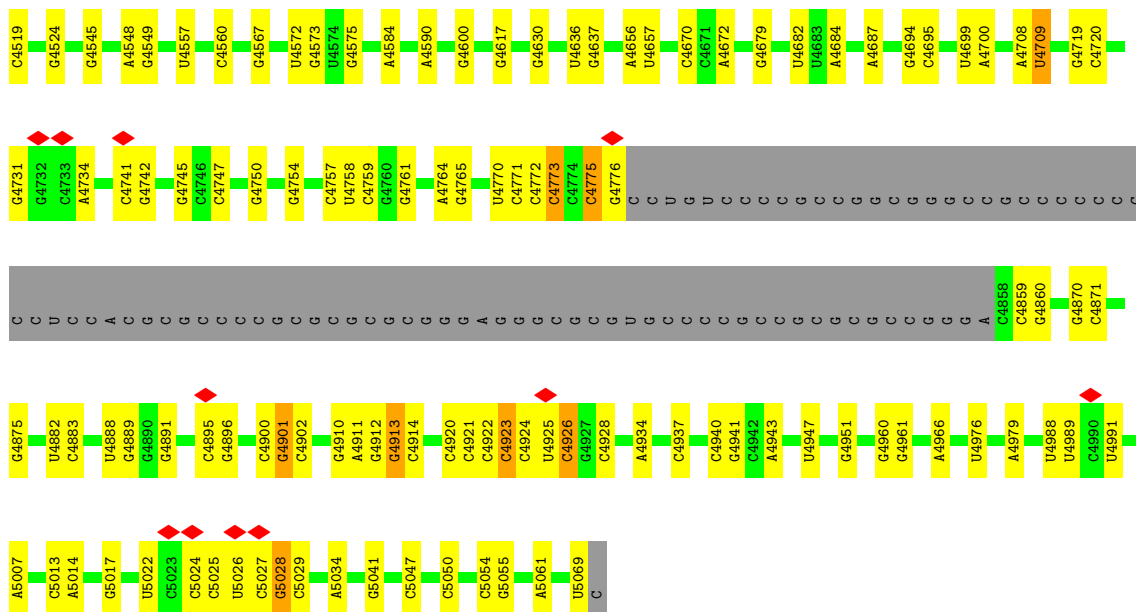
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

- Molecule 1: 28S rRNA

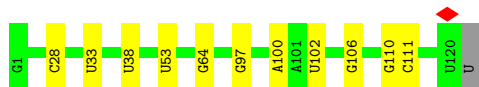
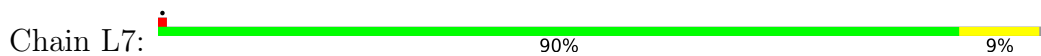




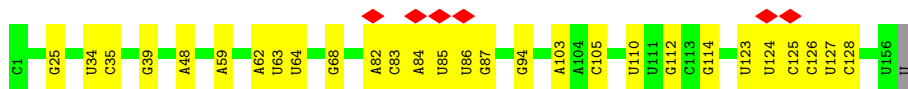
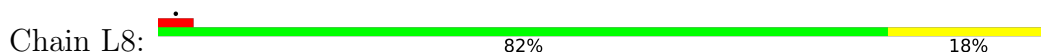
A4273	A4281	G4291	C4303	A4304	G4305	C4314	G4329	G4330	G4331	G4332	C4349	U4354	U4360	G4373	A4376	G4377	A4378	A4379	A4380	C4387	G4391	A4394	C4398	A4422	C4426	G4448	A4449	C4453	A4464	G4475	C4476	A4488	U4500	U4512	A4513	A4268																							
U4120	G4121	G4122	A4127	C4133	C4138	G4139	G4140	G4141	C4142	G4143	G4144	G4145	G4146	G4147	C4148	C4149	G4150	C4162	U4163	G4168	G4169	A4170	G4183	G4191	G4196	G4197	A4203	G4222	G4228	U4229	A4233	C4237	C4241	U4242	A4251	G4254	A4257	U4265																					
C4037	C4038	G4039	C4040	C4041	G4042	G4043	U4044	G4045	A4046	A4047	A4048	U4049	A4050	C4051	C4052	A4053	C4054	U4055	A4056	C4057	U4058	C4059	C4064	G4065	U4066	G4076	G4084	A4085	G4086	G4094	G4095	C4096	G4097	A4098	G4099	C4100	C4101	C4102	C4103	G4104	A4105	G4106	G4107	G4108	G4109	C4110	U4111	C4112	U4113	C4114	G4115	U4116	U4117	C4119					
C3977	C3978	C3979	G3980	G3981	C3982	G3983	C3984	C3985	C3986	C3987	C3988	C3989	C3990	G3991	C3992	U3993	G3994	U3995	C3996	C3997	U3998	C3999	G4000	C4001	C4002	A4003	G4004	G4005	G4006	G4007	C4008	C4009	C4010	G4011	G4012	G4013	G4014	C4015	G4016	G4017	G4018	G4019	U4020	C4021	C4022	G4023	C4024	C4025	G4026	G4027	C4028	C4029	C4030	U4031	G4032	C4033	G4034	G4035	G4036
C3878	G3879	C3882	G3885	G3886	C3887	A3890	A3891	U3892	G3897	A3901	A3906	G3907	A3908	C3909	U3915	C3926	G3938	G3939	G3942	A3943	A3947	U3950	G3951	A3954	G3955	G3956	U3957	G3958	U3959	A3960	G3961	A3962	A3963	U3964	A3965	A3966	G3967	U3968	G3969	G3970	G3971	A3972	G3973	G3974	C3975	C3976													
G3674	U3693	G3710	G3714	U3715	A3727	A3728	U3729	G3735	A3736	C3741	G3750	G3753	G3757	U3758	A3759	A3760	C3761	U3762	C3771	U3772	G3776	U3778	A3784	A3785	U3786	U3802	G3811	C3812	A3813	U3814	A3817	U3818	G3819	G3823	U3838	G3839	U3840	A3887	A3877																				
U3584	G3585	C3587	G3588	G3590	C3591	G3592	C3593	C3594	U3595	A3596	G3597	A3604	C3605	U3606	G3614	G3615	U3616	C3617	C3618	G3626	A3630	A3635	C3636	U3644	U3645	A3646	A3647	A3648	A3662	C3673																													



• Molecule 2: 5S rRNA



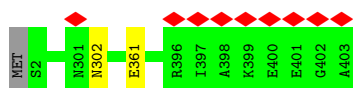
• Molecule 3: 5.8S rRNA



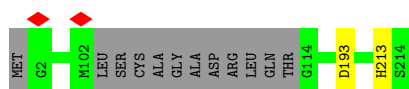
• Molecule 4: 60S ribosomal protein L8



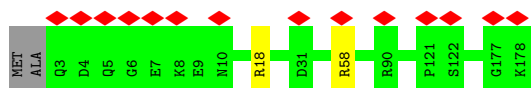
• Molecule 5: 60S ribosomal protein L3



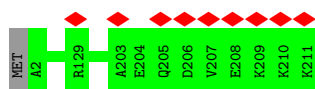
• Molecule 6: 60S ribosomal protein L4



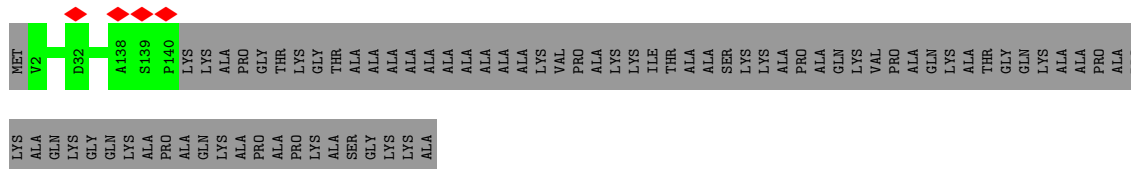
- Molecule 13: 60S ribosomal protein L11



- Molecule 14: 60S ribosomal protein L13



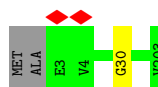
- Molecule 15: 60S ribosomal protein L14



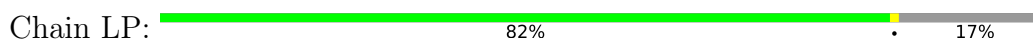
- Molecule 16: 60S ribosomal protein L15

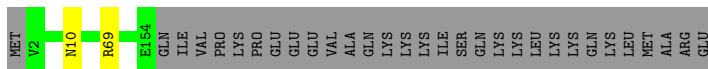


- Molecule 17: 60S ribosomal protein L13a



- Molecule 18: 60S ribosomal protein L17

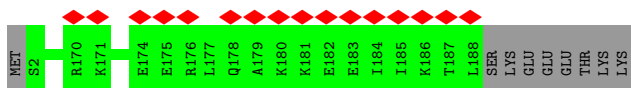




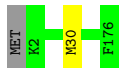
- Molecule 19: 60S ribosomal protein L18



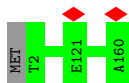
- Molecule 20: 60S ribosomal protein L19



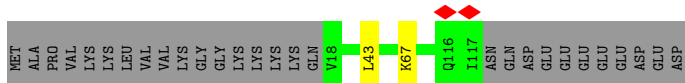
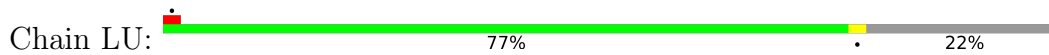
- Molecule 21: 60S ribosomal protein L18a



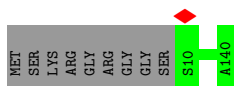
- Molecule 22: 60S ribosomal protein L21



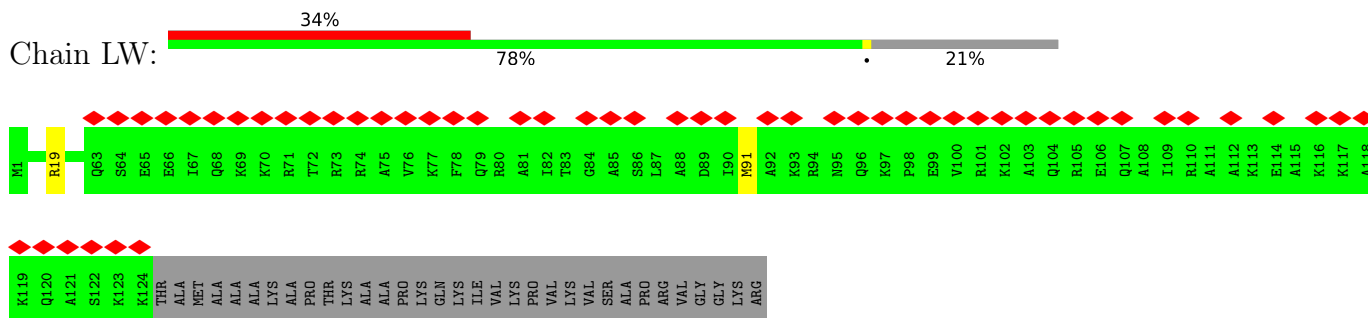
- Molecule 23: 60S ribosomal protein L22



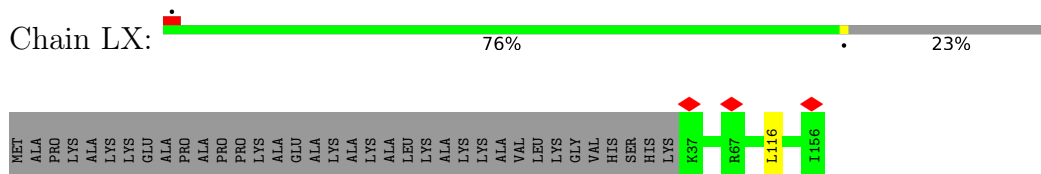
- Molecule 24: 60S ribosomal protein L23



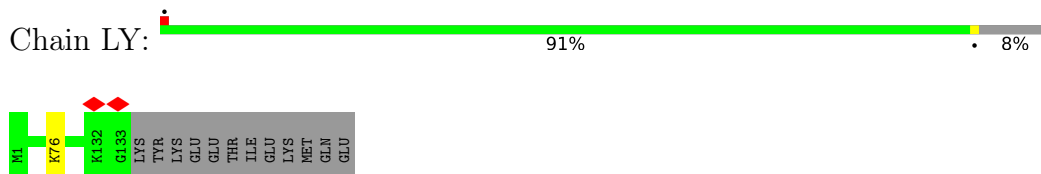
- Molecule 25: 60S ribosomal protein L24



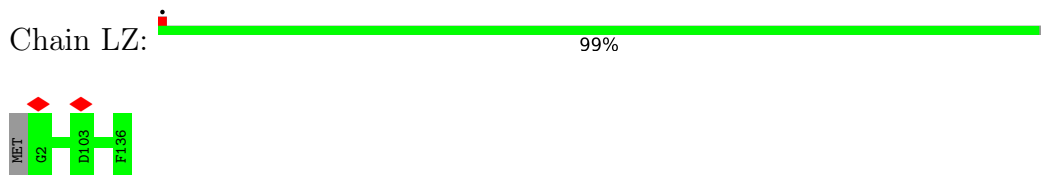
- Molecule 26: 60S ribosomal protein L23a



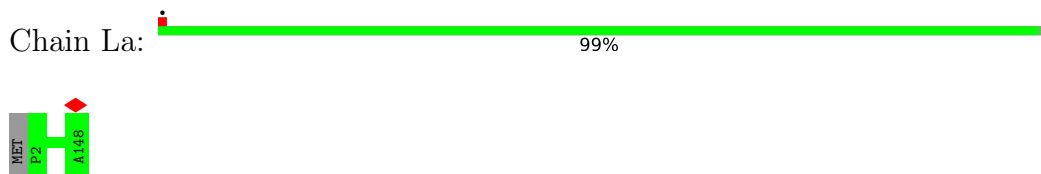
- Molecule 27: 60S ribosomal protein L26



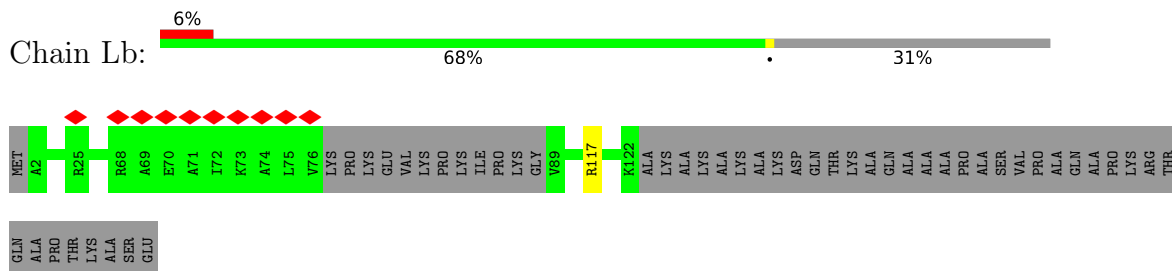
- Molecule 28: 60S ribosomal protein L27



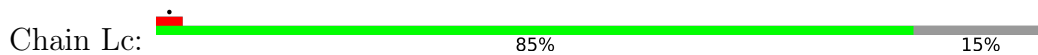
- Molecule 29: 60S ribosomal protein L27a



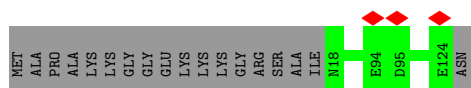
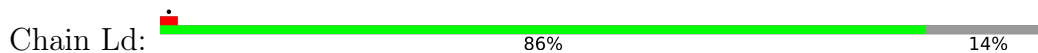
- Molecule 30: 60S ribosomal protein L29



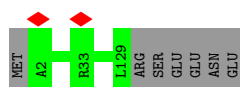
- Molecule 31: 60S ribosomal protein L30



- Molecule 32: 60S ribosomal protein L31



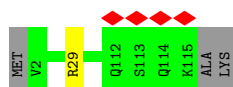
- Molecule 33: 60S ribosomal protein L32



- Molecule 34: 60S ribosomal protein L35a



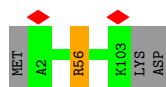
- Molecule 35: 60S ribosomal protein L34



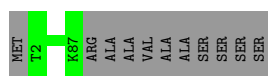
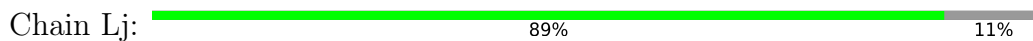
- Molecule 36: 60S ribosomal protein L35



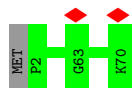
- Molecule 37: 60S ribosomal protein L36



• Molecule 38: 60S ribosomal protein L37



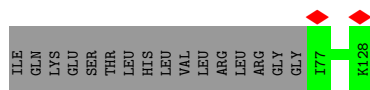
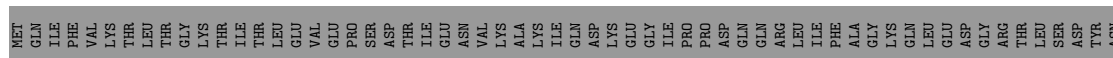
• Molecule 39: 60S ribosomal protein L38



• Molecule 40: 60S ribosomal protein L39



• Molecule 41: Ubiquitin-60S ribosomal protein L40



• Molecule 42: 60S ribosomal protein L41

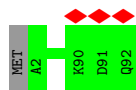


• Molecule 43: 60S ribosomal protein L36a

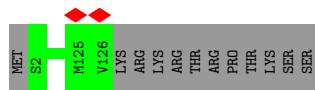


• Molecule 44: 60S ribosomal protein L37a

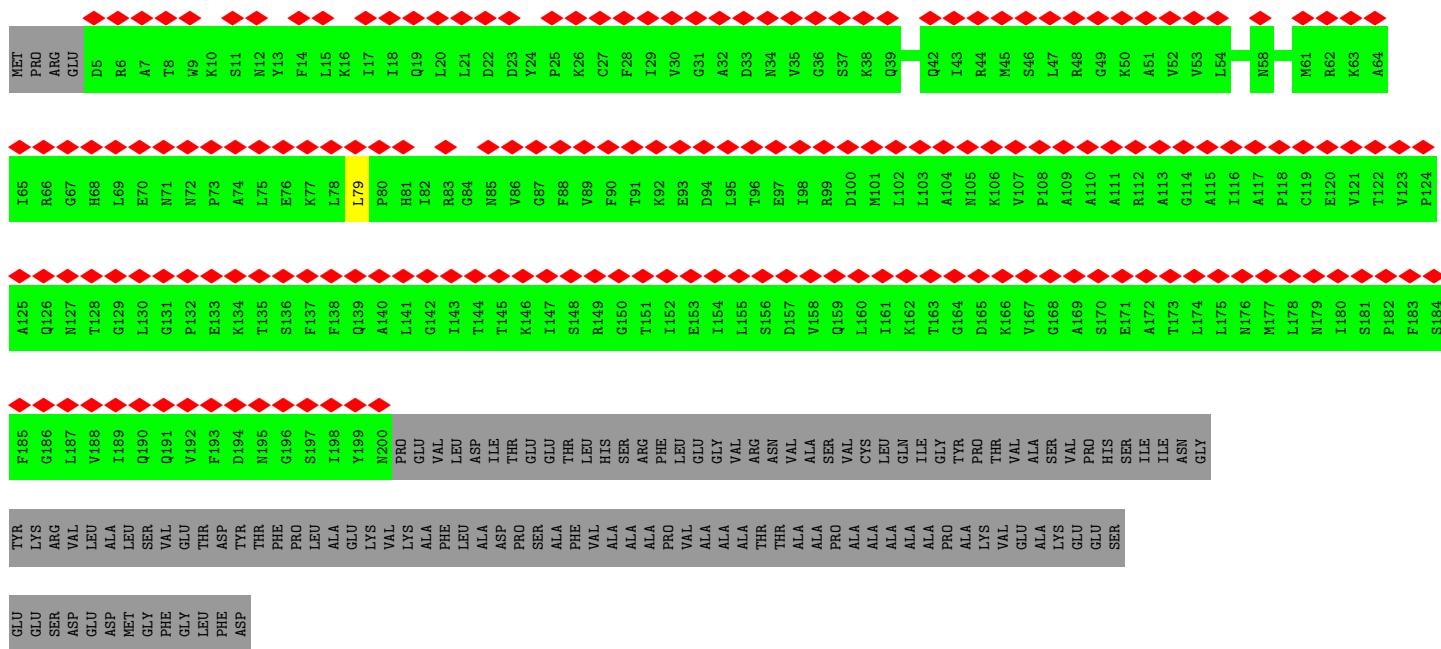




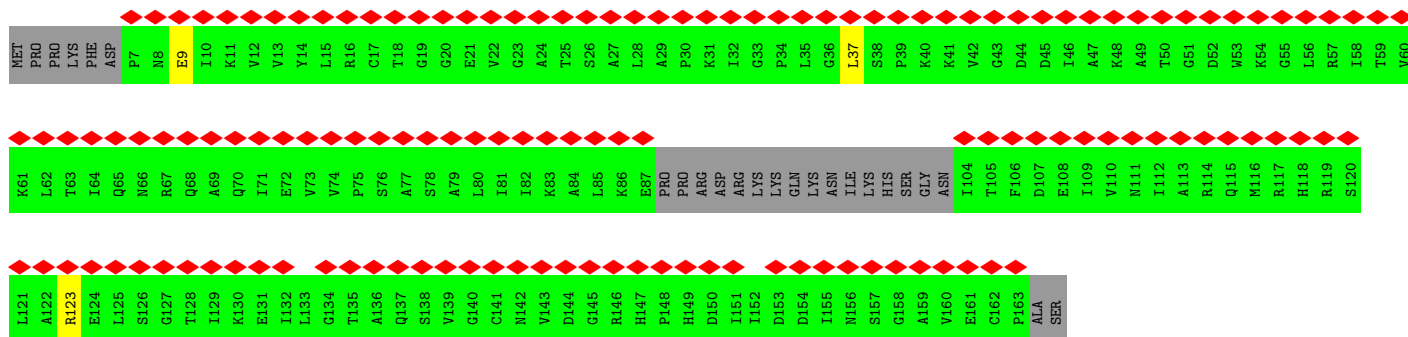
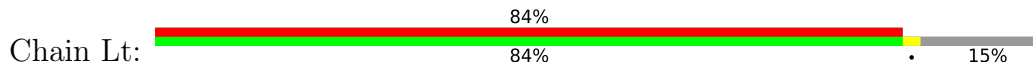
• Molecule 45: 60S ribosomal protein L28



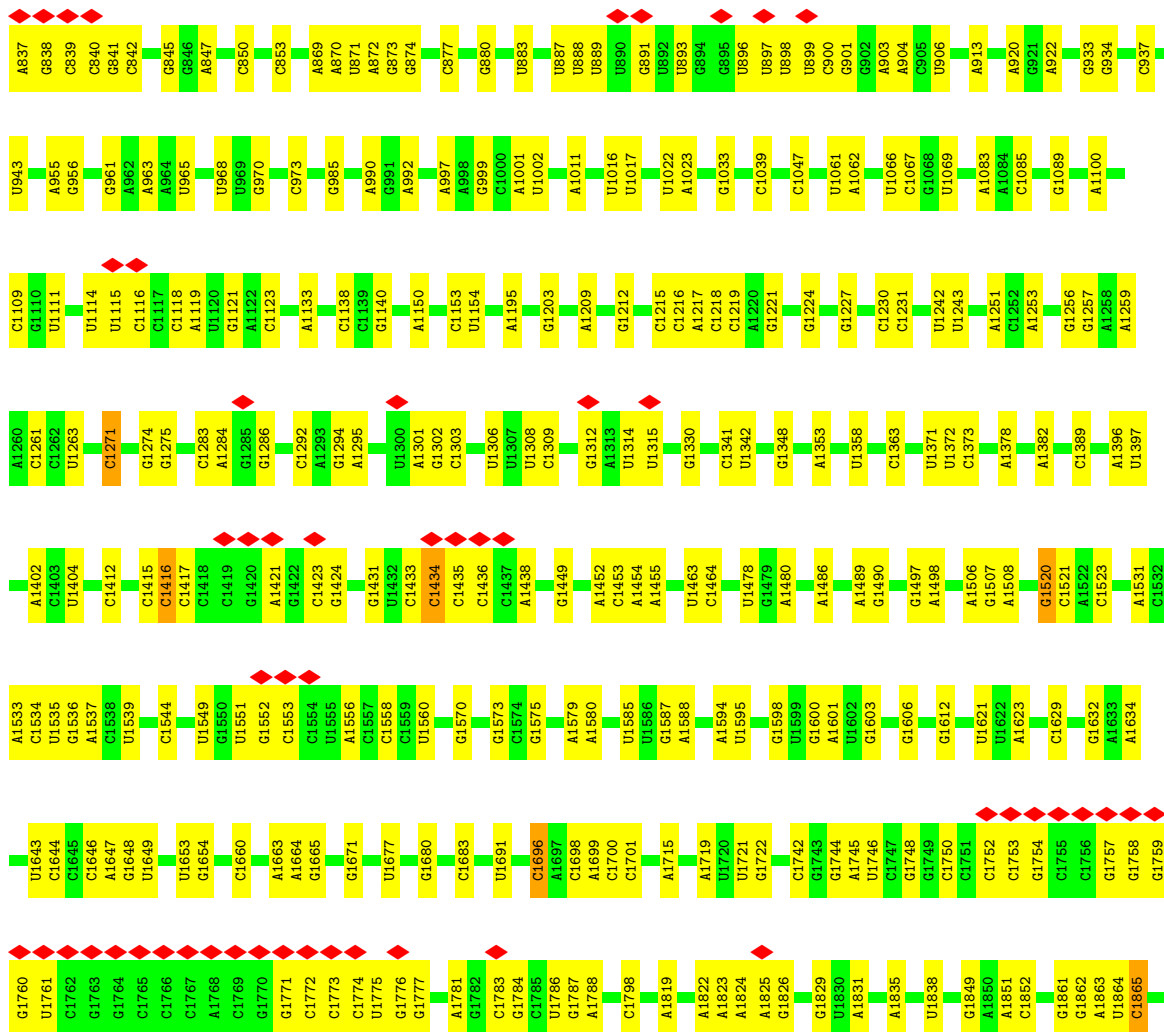
• Molecule 46: Large ribosomal subunit protein uL10



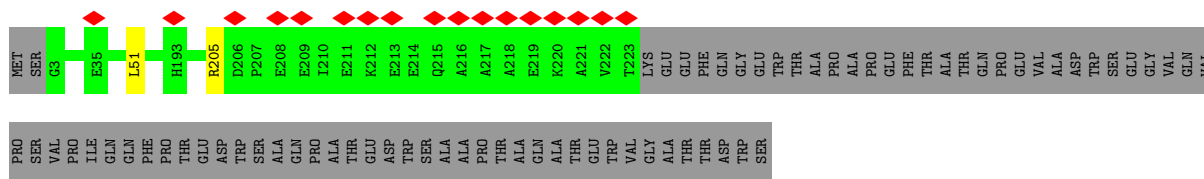
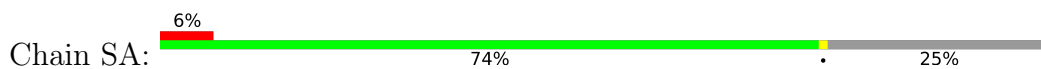
• Molecule 47: 60S ribosomal protein L12



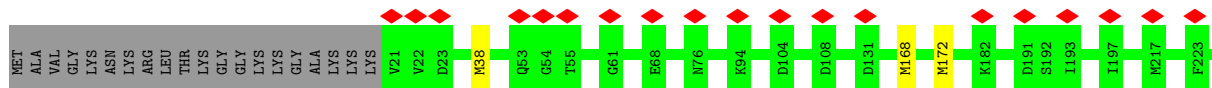
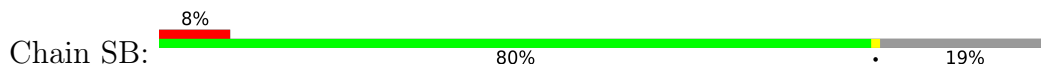
• Molecule 48: 60S ribosomal protein L10a

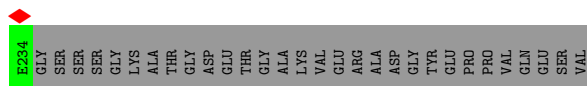


• Molecule 50: 40S ribosomal protein SA

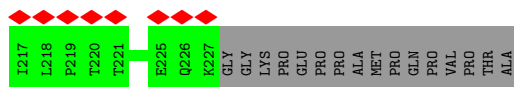
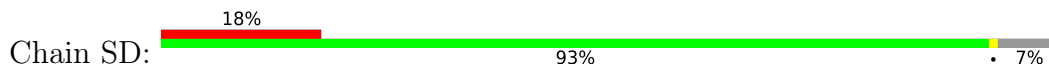


• Molecule 51: 40S ribosomal protein S3a

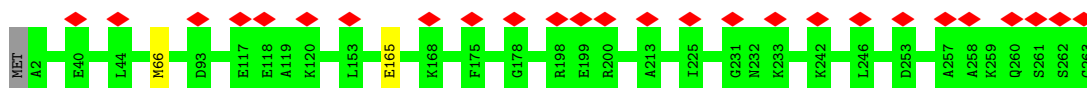




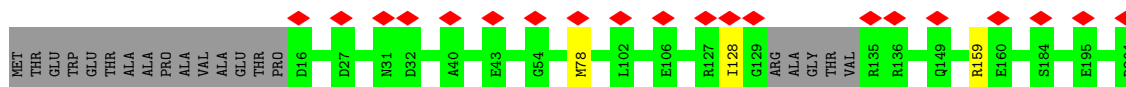
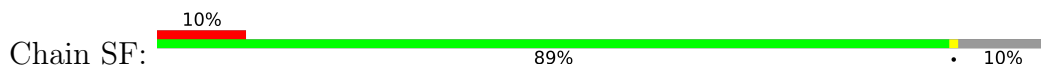
- Molecule 52: 40S ribosomal protein S3



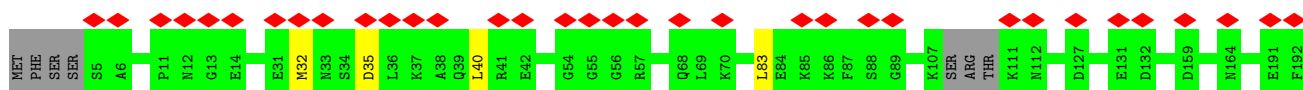
- Molecule 53: 40S ribosomal protein S4, X isoform



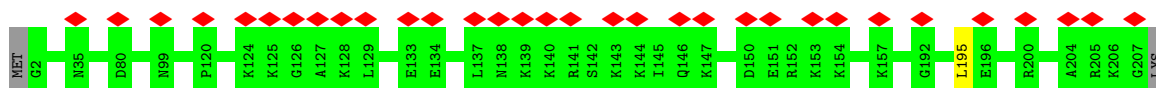
- Molecule 54: 40S ribosomal protein S5



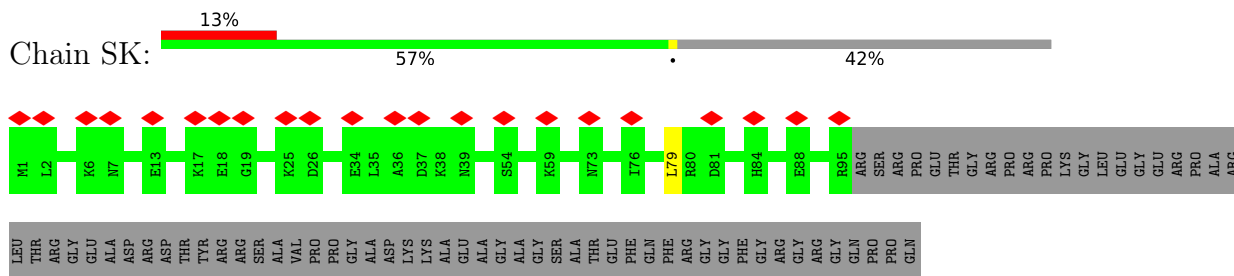
- Molecule 55: 40S ribosomal protein S7



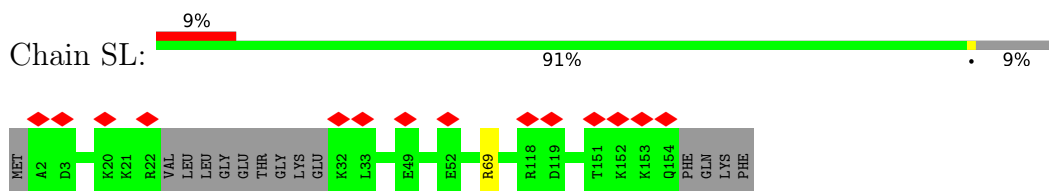
- Molecule 56: 40S ribosomal protein S8



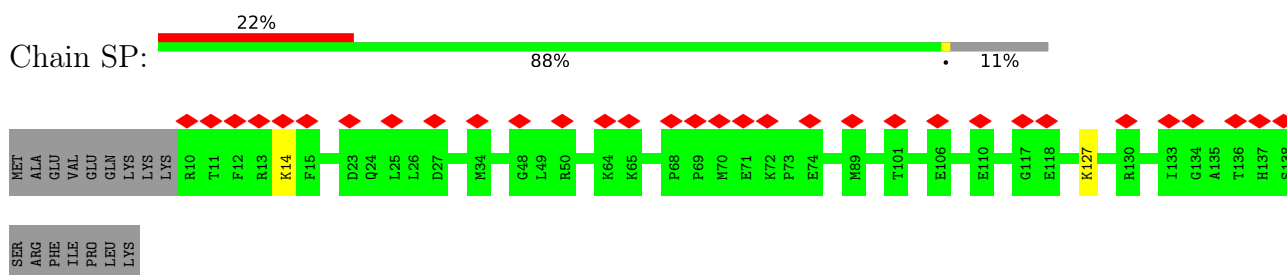
- Molecule 57: 40S ribosomal protein S10



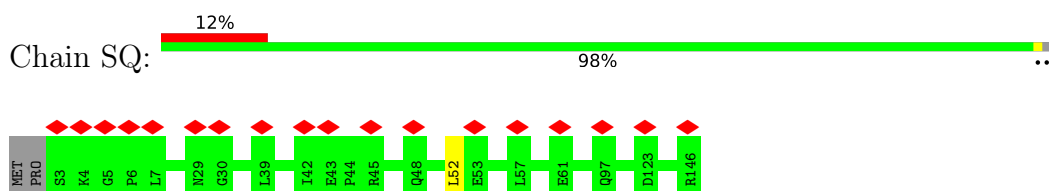
• Molecule 58: 40S ribosomal protein S11



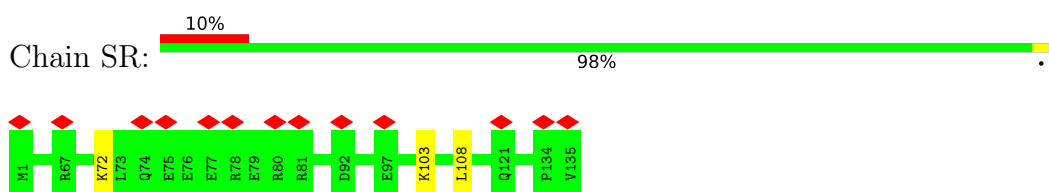
• Molecule 59: 40S ribosomal protein S15



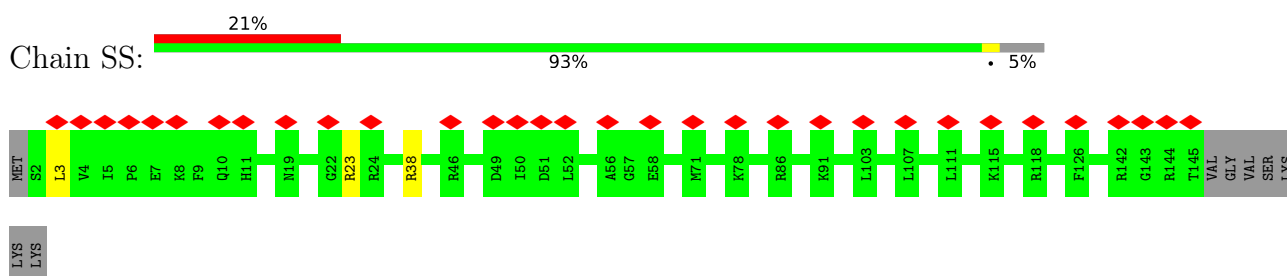
• Molecule 60: 40S ribosomal protein S16



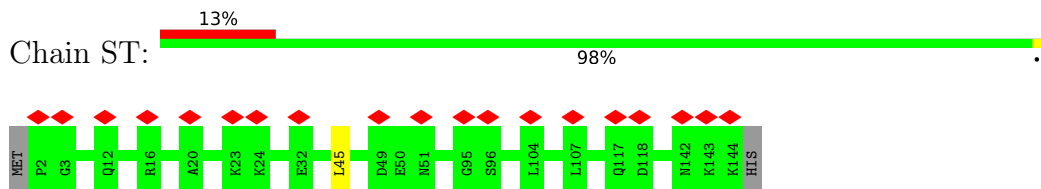
• Molecule 61: 40S ribosomal protein S17



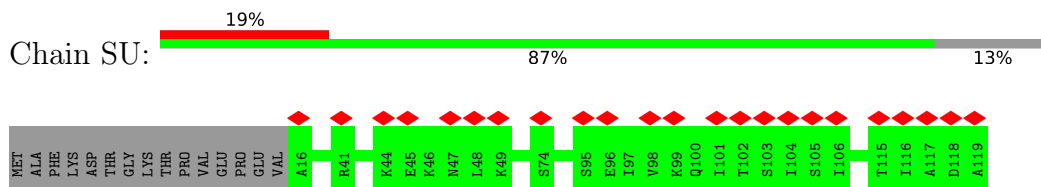
• Molecule 62: 40S ribosomal protein S18



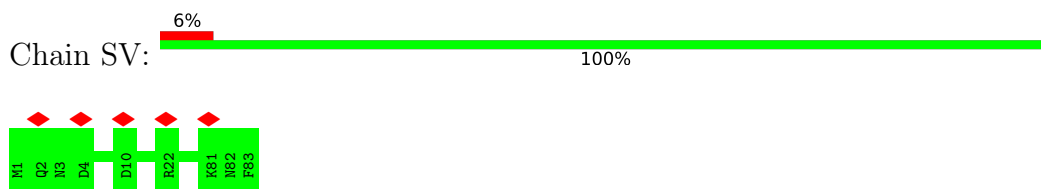
- Molecule 63: 40S ribosomal protein S19



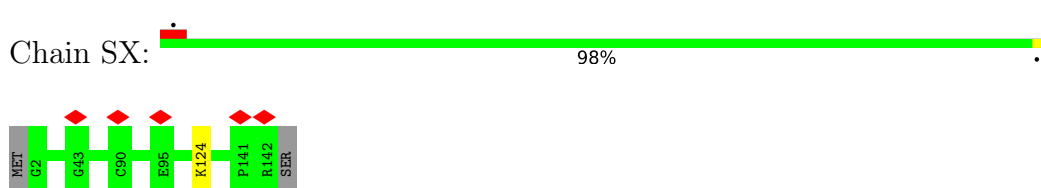
- Molecule 64: 40S ribosomal protein S20



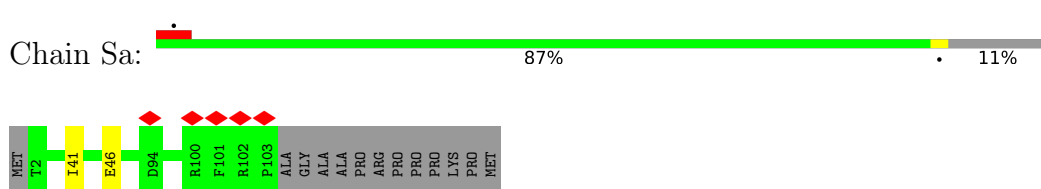
- Molecule 65: 40S ribosomal protein S21



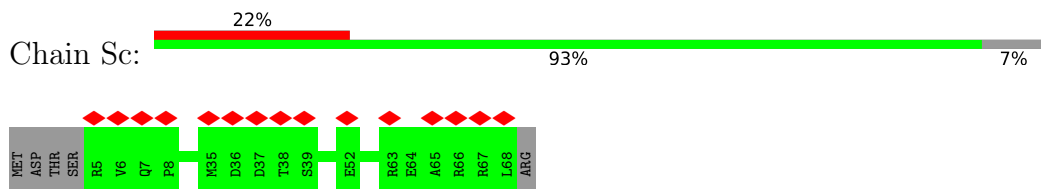
- Molecule 66: 40S ribosomal protein S23



- Molecule 67: 40S ribosomal protein S26

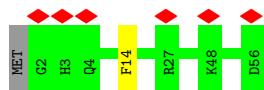


- Molecule 68: 40S ribosomal protein S28

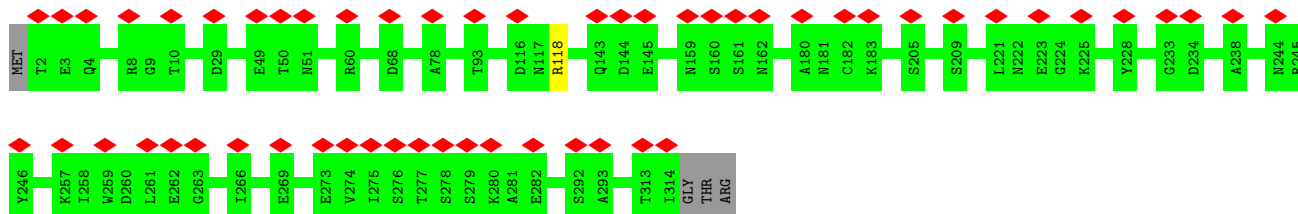


- Molecule 69: 40S ribosomal protein S29

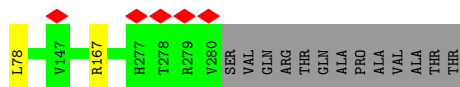
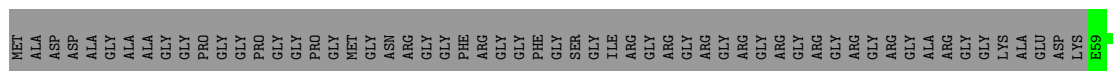
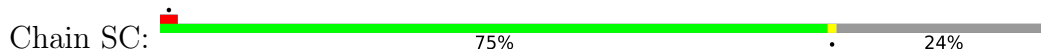




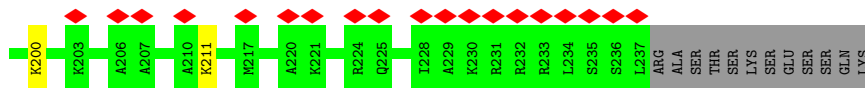
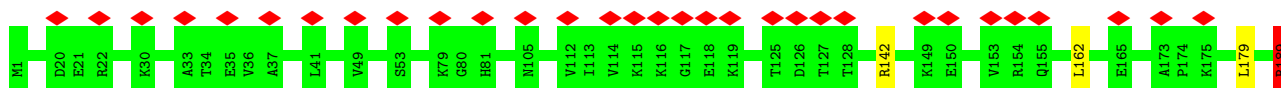
- Molecule 70: Receptor of activated protein C kinase 1



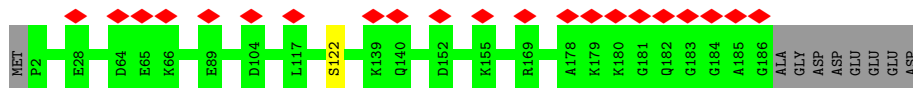
- Molecule 71: 40S ribosomal protein S2



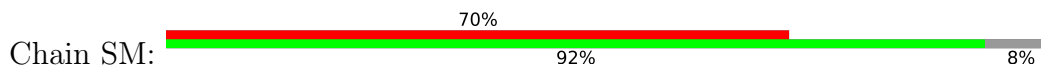
- Molecule 72: 40S ribosomal protein S6

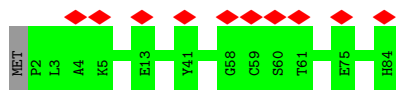


- Molecule 73: 40S ribosomal protein S9



- Molecule 74: 40S ribosomal protein S12

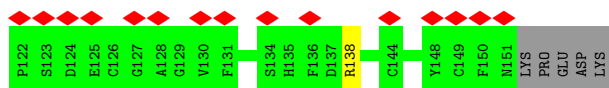
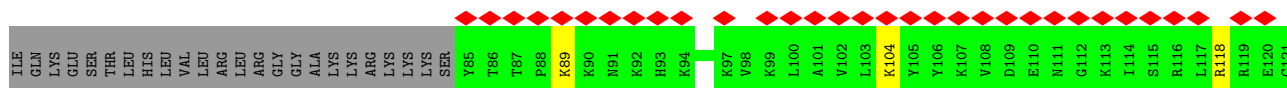
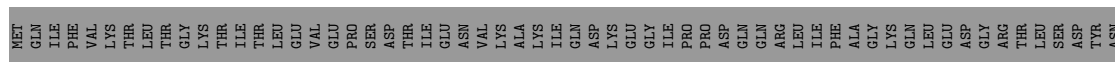




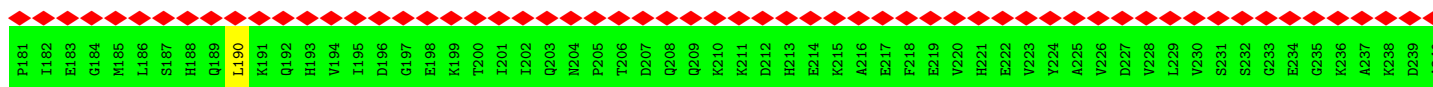
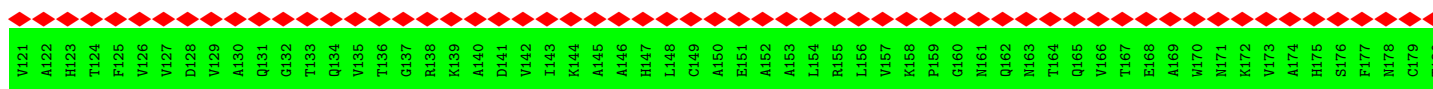
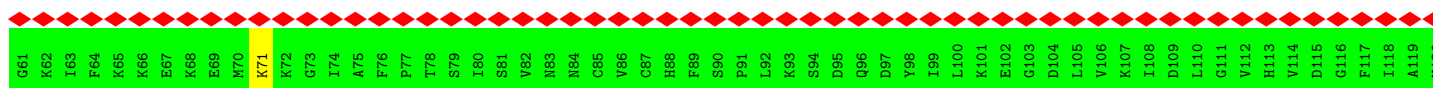
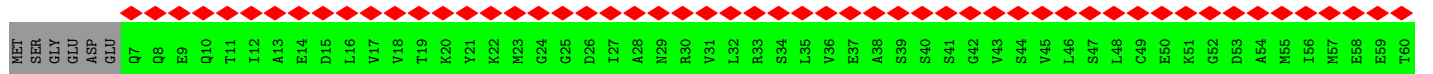
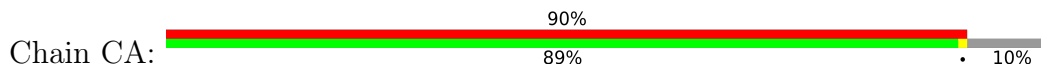
- Molecule 81: 40S ribosomal protein S30

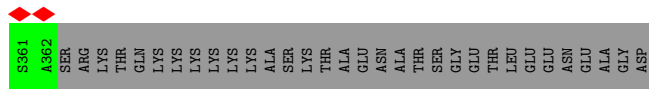
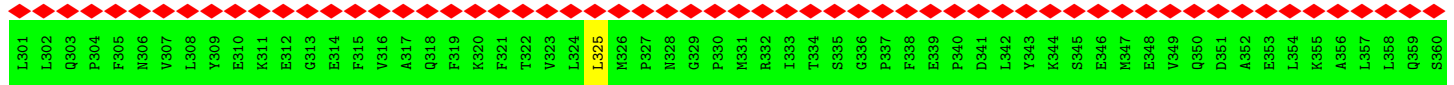


- Molecule 82: Ubiquitin-40S ribosomal protein S27a

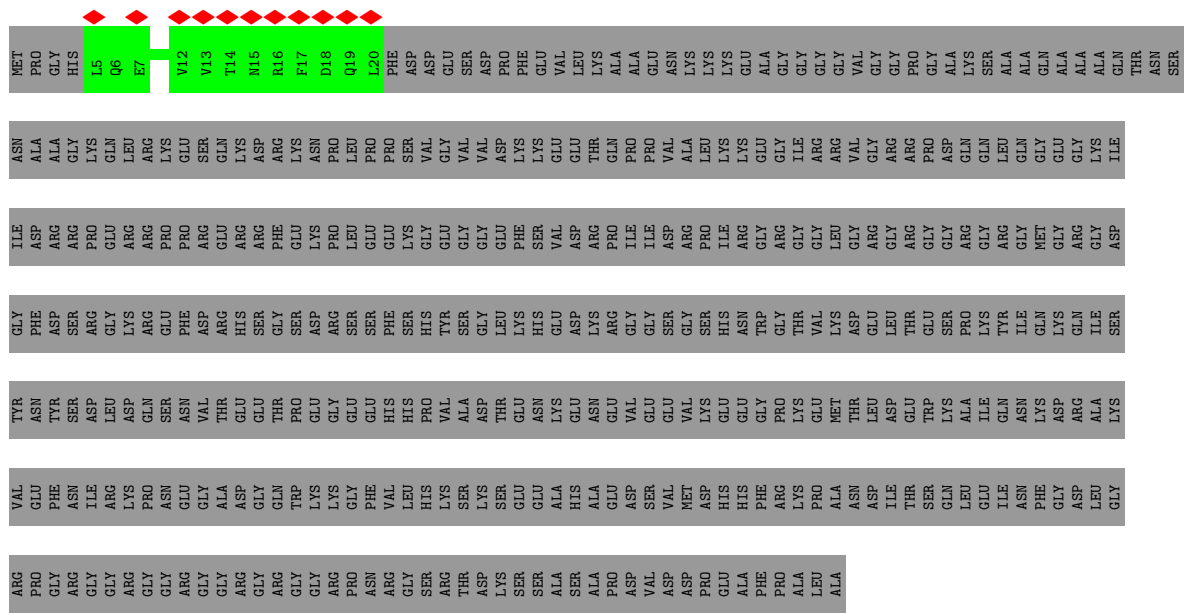


- Molecule 83: Proliferation-associated protein 2G4

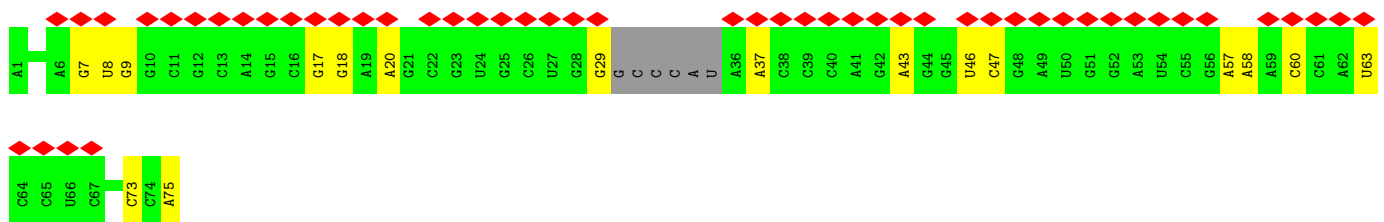




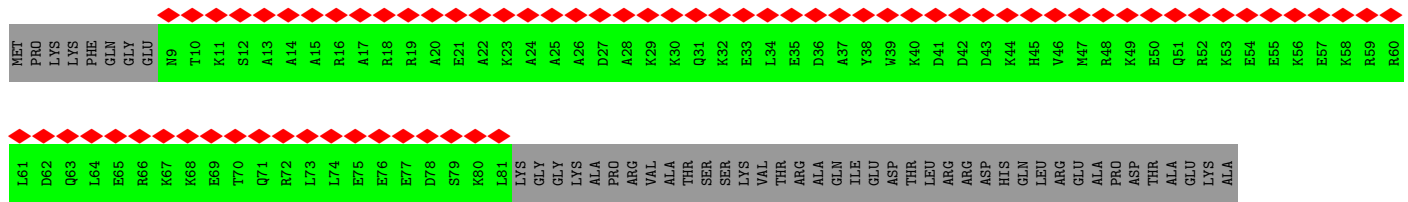
• Molecule 84: SERPINE1 mRNA-binding protein 1



• Molecule 85: E-tRNA



• Molecule 86: Coiled-coil domain-containing protein 124



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

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	46089	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; Relion	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.059	Depositor
Minimum map value	-0.023	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	447.36, 447.36, 447.36	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.932, 0.932, 0.932	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZN, T1C, MLZ

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	L5	0.35	0/89548	0.97	224/139613 (0.2%)
2	L7	0.33	0/2861	0.94	3/4459 (0.1%)
3	L8	0.32	0/3701	0.89	5/5766 (0.1%)
4	LA	0.32	0/1936	0.66	0/2596
5	LB	0.31	0/3306	0.61	0/4424
6	LC	0.31	0/2981	0.63	2/4002 (0.0%)
7	LD	0.30	0/2428	0.56	0/3252
8	LE	0.27	0/1808	0.61	0/2425
9	LF	0.30	0/1905	0.59	0/2539
10	LG	0.30	0/1960	0.60	0/2637
11	LH	0.30	0/1537	0.61	0/2066
12	LI	0.31	0/1677	0.62	1/2237 (0.0%)
13	LJ	0.33	0/1433	0.67	0/1915
14	LL	0.30	0/1732	0.62	0/2315
15	LM	0.29	0/1161	0.62	0/1554
16	LN	0.31	0/1746	0.65	1/2338 (0.0%)
17	LO	0.30	0/1682	0.58	0/2250
18	LP	0.28	0/1268	0.56	0/1701
19	LQ	0.28	0/1537	0.63	0/2052
20	LR	0.27	0/1582	0.60	0/2091
21	LS	0.34	0/1493	0.63	1/2003 (0.0%)
22	LT	0.33	0/1326	0.61	0/1770
23	LU	0.31	0/830	0.66	1/1114 (0.1%)
24	LV	0.30	0/993	0.59	0/1332
25	LW	0.34	0/1030	0.69	1/1364 (0.1%)
26	LX	0.28	0/1002	0.59	1/1345 (0.1%)
27	LY	0.31	0/1123	0.61	1/1493 (0.1%)
28	LZ	0.29	0/1130	0.57	0/1507
29	La	0.32	0/1191	0.59	0/1591
30	Lb	0.27	0/895	0.62	0/1182
31	Lc	0.31	0/774	0.57	0/1038
32	Ld	0.30	0/903	0.64	0/1216

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	Le	0.29	0/1071	0.60	0/1429
34	Lf	0.31	0/895	0.63	0/1198
35	Lg	0.29	0/916	0.61	0/1220
36	Lh	0.27	0/1023	0.60	0/1351
37	Li	0.28	0/843	0.63	0/1115
38	Lj	0.29	0/720	0.64	0/952
39	Lk	0.31	0/575	0.58	0/761
40	Ll	0.26	0/454	0.57	0/599
41	Lm	0.27	0/425	0.58	0/561
42	Ln	0.29	0/231	0.73	0/294
43	Lo	0.30	0/876	0.60	0/1156
44	Lp	0.30	0/718	0.57	0/953
45	Lr	0.28	0/1017	0.59	0/1364
46	Ls	0.27	0/1519	0.63	1/2052 (0.0%)
47	Lt	0.28	0/1058	0.72	1/1430 (0.1%)
48	Lz	0.25	0/1017	0.47	0/1416
49	S2	0.31	0/41241	0.98	105/64258 (0.2%)
50	SA	0.31	0/1778	0.62	1/2416 (0.0%)
51	SB	0.33	0/1765	0.69	4/2362 (0.2%)
52	SD	0.35	0/1793	0.69	1/2414 (0.0%)
53	SE	0.29	0/2118	0.63	0/2849
54	SF	0.31	0/1481	0.67	1/1988 (0.1%)
55	SH	0.36	0/1519	0.69	4/2033 (0.2%)
56	SI	0.29	0/1715	0.64	1/2287 (0.0%)
57	SK	0.41	0/823	0.75	1/1111 (0.1%)
58	SL	0.28	0/1202	0.57	0/1606
59	SP	0.35	0/1082	0.76	0/1446
60	SQ	0.30	0/1160	0.71	1/1553 (0.1%)
61	SR	0.36	0/1105	0.75	2/1484 (0.1%)
62	SS	0.36	0/1208	0.77	1/1618 (0.1%)
63	ST	0.39	0/1131	0.66	0/1515
64	SU	0.36	0/831	0.71	0/1115
65	SV	0.27	0/643	0.55	0/860
66	SX	0.30	0/1116	0.62	0/1490
67	Sa	0.30	0/836	0.69	1/1121 (0.1%)
68	Sc	0.30	0/508	0.76	0/680
69	Sd	0.34	0/470	0.71	0/623
70	Sg	0.29	0/2493	0.66	0/3394
71	SC	0.35	0/1762	0.64	1/2381 (0.0%)
72	SG	0.31	0/1946	0.74	4/2590 (0.2%)
73	SJ	0.34	0/1550	0.66	0/2069
74	SM	0.24	0/603	0.40	0/837
75	SN	0.29	0/1232	0.58	0/1656

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
76	SO	0.31	0/1023	0.67	0/1372
77	SW	0.30	0/1051	0.62	0/1406
78	SY	0.29	0/1044	0.61	1/1388 (0.1%)
79	SZ	0.42	0/604	1.19	5/810 (0.6%)
80	Sb	0.28	0/665	0.52	0/891
81	Se	0.27	0/465	0.65	0/612
82	Sf	0.34	0/560	0.74	0/745
83	CA	0.33	0/2810	0.67	4/3780 (0.1%)
84	CB	0.31	0/128	0.62	0/171
85	CC	0.26	0/1658	0.88	0/2582
86	CE	0.26	0/616	0.48	0/812
All	All	0.32	0/239572	0.86	381/351363 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
4	LA	0	1
13	LJ	0	1
17	LO	0	1
25	LW	0	1
30	Lb	0	1
35	Lg	0	1
37	Li	0	1
50	SA	0	1
53	SE	0	1
54	SF	0	1
62	SS	0	2
72	SG	0	1
All	All	0	13

There are no bond length outliers.

All (381) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	485	C	C2-N1-C1'	13.35	133.49	118.80
49	S2	1263	U	C5-C4-O4	-12.91	118.16	125.90
1	L5	1378	C	N1-C2-O2	12.87	126.62	118.90
79	SZ	88	LEU	CB-CG-CD2	-12.78	89.28	111.00
1	L5	1367	C	N1-C2-O2	12.39	126.33	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	1367	C	C2-N1-C1'	11.40	131.34	118.80
1	L5	4923	C	N3-C2-O2	-10.23	114.74	121.90
1	L5	1378	C	C2-N1-C1'	10.21	130.03	118.80
1	L5	1367	C	N3-C2-O2	-10.18	114.77	121.90
1	L5	4149	C	N3-C2-O2	-10.13	114.81	121.90
1	L5	1378	C	N3-C2-O2	-10.06	114.85	121.90
79	SZ	83	LEU	CB-CG-CD2	-10.02	93.96	111.00
1	L5	485	C	C6-N1-C1'	-10.00	108.80	120.80
49	S2	1535	U	N3-C4-O4	-9.82	112.53	119.40
49	S2	1453	C	C2-N1-C1'	9.50	129.25	118.80
1	L5	2409	U	C2-N1-C1'	9.50	129.10	117.70
1	L5	4921	C	N3-C2-O2	-9.09	115.54	121.90
1	L5	4138	C	N3-C2-O2	-9.08	115.54	121.90
49	S2	1453	C	N1-C2-O2	8.90	124.24	118.90
49	S2	331	C	C6-N1-C2	-8.77	116.79	120.30
1	L5	4928	C	C2-N1-C1'	8.61	128.27	118.80
1	L5	515	C	N1-C2-O2	8.56	124.03	118.90
1	L5	1367	C	C6-N1-C2	-8.42	116.93	120.30
51	SB	168	MET	CB-CG-SD	8.35	137.45	112.40
1	L5	485	C	N1-C2-O2	8.29	123.87	118.90
1	L5	3715	U	N3-C2-O2	-8.26	116.42	122.20
51	SB	172	MET	CB-CG-SD	-8.23	87.70	112.40
49	S2	293	C	N1-C2-O2	8.21	123.82	118.90
1	L5	1252	C	N3-C2-O2	-8.19	116.17	121.90
1	L5	417	G	O4'-C1'-N9	8.17	114.74	108.20
49	S2	293	C	C2-N1-C1'	8.17	127.79	118.80
1	L5	2409	U	N1-C2-O2	8.12	128.48	122.80
1	L5	1241	C	N1-C2-O2	8.09	123.75	118.90
1	L5	4923	C	N1-C2-O2	8.07	123.74	118.90
1	L5	2409	U	N3-C2-O2	-7.90	116.67	122.20
1	L5	1241	C	C2-N1-C1'	7.89	127.48	118.80
49	S2	1309	C	C2-N1-C1'	7.68	127.24	118.80
1	L5	2627	C	C2-N1-C1'	7.67	127.24	118.80
1	L5	1216	C	C2-N1-C1'	7.66	127.23	118.80
1	L5	2814	C	N1-C2-O2	7.65	123.49	118.90
1	L5	1367	C	C6-N1-C1'	-7.64	111.63	120.80
79	SZ	44	LEU	CA-CB-CG	7.62	132.82	115.30
1	L5	4303	C	C2-N1-C1'	7.45	127.00	118.80
49	S2	1520	G	C4-N9-C1'	7.43	136.16	126.50
1	L5	3715	U	C2-N1-C1'	7.40	126.58	117.70
1	L5	925	C	N3-C2-O2	-7.37	116.74	121.90
3	L8	128	C	C2-N1-C1'	7.36	126.89	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	126	G	N1-C6-O6	-7.29	115.53	119.90
1	L5	515	C	C2-N1-C1'	7.19	126.71	118.80
1	L5	1378	C	C6-N1-C2	-7.19	117.43	120.30
1	L5	4758	U	C2-N1-C1'	7.17	126.30	117.70
1	L5	2410	C	C2-N1-C1'	7.16	126.67	118.80
72	SG	189	ARG	CB-CG-CD	-7.13	93.07	111.60
1	L5	4924	C	N3-C2-O2	-7.11	116.92	121.90
1	L5	3741	C	N1-C2-O2	7.09	123.16	118.90
49	S2	1047	C	C5-C6-N1	7.09	124.54	121.00
1	L5	2018	C	C5-C6-N1	7.05	124.53	121.00
83	CA	257	LEU	CB-CG-CD2	-7.03	99.05	111.00
49	S2	583	C	N1-C2-O2	7.02	123.11	118.90
1	L5	3741	C	N3-C2-O2	-6.97	117.02	121.90
49	S2	1271	C	N1-C2-O2	6.96	123.08	118.90
1	L5	4229	U	N3-C2-O2	-6.96	117.33	122.20
49	S2	1535	U	C5-C4-O4	6.95	130.07	125.90
1	L5	925	C	N1-C2-O2	6.90	123.04	118.90
1	L5	3715	U	N1-C2-O2	6.89	127.63	122.80
1	L5	1378	C	C6-N1-C1'	-6.89	112.53	120.80
49	S2	688	U	P-O3'-C3'	6.89	127.97	119.70
49	S2	1389	C	C2-N1-C1'	6.89	126.38	118.80
49	S2	180	G	C2-N3-C4	6.87	115.33	111.90
1	L5	2814	C	C2-N1-C1'	6.86	126.35	118.80
1	L5	2494	U	N1-C2-O2	6.86	127.60	122.80
49	S2	1520	G	N3-C4-N9	6.86	130.11	126.00
49	S2	1520	G	C8-N9-C1'	-6.82	118.13	127.00
1	L5	4928	C	N1-C2-O2	6.80	122.98	118.90
55	SH	83	LEU	CB-CG-CD2	6.79	122.54	111.00
1	L5	1216	C	N1-C2-O2	6.78	122.97	118.90
1	L5	1191	C	N3-C2-O2	-6.76	117.17	121.90
1	L5	1241	C	N3-C2-O2	-6.76	117.17	121.90
1	L5	2627	C	N1-C2-O2	6.76	122.95	118.90
67	Sa	41	ILE	CG1-CB-CG2	-6.74	96.57	111.40
1	L5	485	C	C5-C6-N1	6.72	124.36	121.00
49	S2	1536	G	C2-N3-C4	6.72	115.26	111.90
1	L5	4107	G	N3-C4-C5	-6.71	125.24	128.60
54	SF	159	ARG	CG-CD-NE	6.71	125.90	111.80
1	L5	963	G	C4-N9-C1'	6.70	135.22	126.50
49	S2	1453	C	N3-C2-O2	-6.70	117.21	121.90
1	L5	4758	U	N1-C2-O2	6.62	127.43	122.80
49	S2	1047	C	C2-N1-C1'	6.61	126.07	118.80
1	L5	4149	C	C6-N1-C2	-6.60	117.66	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	293	C	N3-C2-O2	-6.60	117.28	121.90
49	S2	1453	C	C6-N1-C1'	-6.57	112.92	120.80
1	L5	485	C	C6-N1-C2	-6.57	117.67	120.30
12	LI	193	ASP	CB-CG-OD2	-6.55	112.40	118.30
1	L5	4572	U	C2-N1-C1'	6.54	125.55	117.70
1	L5	4107	G	C4-N9-C1'	6.54	135.00	126.50
49	S2	1416	C	C5-C6-N1	6.53	124.27	121.00
1	L5	209	U	C2-N1-C1'	6.51	125.52	117.70
1	L5	1193	C	C2-N1-C1'	6.51	125.96	118.80
49	S2	1022	U	C2-N1-C1'	6.49	125.49	117.70
1	L5	4303	C	N1-C2-O2	6.48	122.79	118.90
49	S2	168	C	N1-C2-O2	6.48	122.79	118.90
1	L5	4557	U	N3-C2-O2	-6.45	117.69	122.20
1	L5	4928	C	N3-C2-O2	-6.44	117.39	121.90
1	L5	4149	C	N1-C2-O2	6.43	122.76	118.90
1	L5	2494	U	N3-C2-O2	-6.41	117.72	122.20
83	CA	257	LEU	CB-CG-CD1	-6.38	100.16	111.00
1	L5	1367	C	C5-C6-N1	6.34	124.17	121.00
1	L5	4557	U	N1-C2-O2	6.34	127.24	122.80
49	S2	1261	C	N1-C2-O2	6.34	122.70	118.90
1	L5	2708	U	C2-N1-C1'	6.33	125.30	117.70
3	L8	64	U	N3-C2-O2	-6.31	117.78	122.20
1	L5	4926	C	C2-N1-C1'	6.28	125.71	118.80
1	L5	2262	G	C4-N9-C1'	6.28	134.67	126.50
1	L5	2760	G	P-O3'-C3'	6.28	127.24	119.70
1	L5	1762	C	C2-N1-C1'	6.27	125.69	118.80
1	L5	4303	C	N3-C2-O2	-6.25	117.53	121.90
1	L5	2708	U	N1-C2-O2	6.24	127.17	122.80
1	L5	220	C	C2-N1-C1'	6.24	125.66	118.80
1	L5	4557	U	C2-N1-C1'	6.23	125.17	117.70
6	LC	319	LEU	CA-CB-CG	6.23	129.62	115.30
49	S2	459	C	C2-N1-C1'	6.22	125.64	118.80
1	L5	2528	G	C4-N9-C1'	6.22	134.58	126.50
1	L5	673	C	C2-N1-C1'	6.21	125.63	118.80
61	SR	108	LEU	CA-CB-CG	6.21	129.58	115.30
1	L5	4928	C	C6-N1-C2	-6.21	117.82	120.30
79	SZ	88	LEU	CA-CB-CG	6.20	129.55	115.30
1	L5	4926	C	N1-C2-O2	6.19	122.61	118.90
1	L5	963	G	N3-C4-N9	6.19	129.71	126.00
1	L5	2409	U	C6-N1-C1'	-6.18	112.55	121.20
47	Lt	37	LEU	CA-CB-CG	6.18	129.50	115.30
1	L5	4758	U	N3-C2-O2	-6.17	117.88	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1453	C	C6-N1-C2	-6.17	117.83	120.30
1	L5	914	U	P-O3'-C3'	6.15	127.08	119.70
49	S2	1416	C	C6-N1-C2	-6.15	117.84	120.30
49	S2	1263	U	N3-C4-C5	6.14	118.29	114.60
1	L5	1216	C	N3-C2-O2	-6.14	117.60	121.90
1	L5	2814	C	N3-C2-O2	-6.14	117.61	121.90
1	L5	1182	C	N1-C2-O2	6.11	122.57	118.90
1	L5	453	G	C4-N9-C1'	6.11	134.44	126.50
49	S2	1219	C	N1-C2-O2	6.11	122.56	118.90
1	L5	1182	C	C2-N1-C1'	6.10	125.51	118.80
1	L5	4926	C	O4'-C1'-N1	6.10	113.08	108.20
1	L5	4303	C	C6-N1-C2	-6.09	117.86	120.30
1	L5	1762	C	N1-C2-O2	6.08	122.55	118.90
52	SD	142	LEU	CB-CG-CD2	6.07	121.31	111.00
46	Ls	79	LEU	CA-CB-CG	6.06	129.24	115.30
49	S2	537	C	C2-N1-C1'	6.05	125.45	118.80
1	L5	1378	C	C5-C6-N1	6.05	124.02	121.00
49	S2	126	G	C5-C6-O6	6.04	132.22	128.60
1	L5	489	C	C2-N1-C1'	6.02	125.42	118.80
49	S2	1309	C	N1-C2-O2	6.00	122.50	118.90
49	S2	1453	C	C5-C6-N1	5.99	124.00	121.00
1	L5	4709	U	N1-C2-O2	5.99	126.99	122.80
1	L5	77	U	N3-C2-O2	-5.98	118.01	122.20
1	L5	4773	C	N1-C2-O2	5.97	122.48	118.90
1	L5	472	C	C2-N1-C1'	5.96	125.36	118.80
49	S2	1315	U	C2-N1-C1'	5.96	124.86	117.70
49	S2	1536	G	N9-C4-C5	5.96	107.78	105.40
1	L5	504	G	C2-N3-C4	5.94	114.87	111.90
1	L5	453	G	N3-C4-C5	-5.93	125.64	128.60
1	L5	1216	C	C6-N1-C2	-5.93	117.93	120.30
1	L5	2255	C	C2-N1-C1'	5.92	125.32	118.80
49	S2	1263	U	N3-C4-O4	5.92	123.54	119.40
1	L5	963	G	C8-N9-C1'	-5.91	119.32	127.00
1	L5	4926	C	C6-N1-C2	-5.89	117.94	120.30
1	L5	2675	G	P-O3'-C3'	5.89	126.76	119.70
1	L5	2820	C	N1-C2-O2	5.88	122.42	118.90
16	LN	134	LEU	CA-CB-CG	5.87	128.81	115.30
27	LY	76	LYS	CD-CE-NZ	-5.87	98.20	111.70
1	L5	2627	C	C6-N1-C2	-5.85	117.96	120.30
1	L5	2262	G	N3-C4-N9	5.85	129.51	126.00
49	S2	427	U	C2-N1-C1'	5.84	124.71	117.70
1	L5	1726	U	N3-C2-O2	-5.84	118.11	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4138	C	N1-C2-O2	5.84	122.41	118.90
3	L8	128	C	C5-C6-N1	5.84	123.92	121.00
49	S2	1315	U	N1-C2-O2	5.83	126.88	122.80
1	L5	753	C	C2-N1-C1'	5.83	125.22	118.80
1	L5	3909	C	C6-N1-C2	-5.82	117.97	120.30
1	L5	1251	C	N1-C2-O2	5.81	122.39	118.90
1	L5	115	C	C2-N1-C1'	5.80	125.18	118.80
1	L5	4107	G	N1-C6-O6	-5.80	116.42	119.90
49	S2	1261	C	N3-C2-O2	-5.80	117.84	121.90
1	L5	963	G	N3-C4-C5	-5.80	125.70	128.60
1	L5	4476	C	C2-N1-C1'	5.80	125.17	118.80
49	S2	1123	C	C2-N1-C1'	5.79	125.17	118.80
1	L5	2262	G	N3-C4-C5	-5.79	125.70	128.60
1	L5	4229	U	N1-C2-O2	5.79	126.85	122.80
49	S2	1219	C	C6-N1-C2	-5.79	117.98	120.30
49	S2	293	C	C6-N1-C2	-5.78	117.99	120.30
1	L5	3772	U	C2-N1-C1'	5.77	124.62	117.70
61	SR	103	LYS	CD-CE-NZ	-5.76	98.44	111.70
2	L7	102	U	N1-C2-O2	5.76	126.83	122.80
49	S2	1263	U	O5'-P-OP2	5.76	117.61	110.70
1	L5	112	C	C2-N1-C1'	5.76	125.13	118.80
1	L5	2033	A	P-O3'-C3'	5.76	126.61	119.70
1	L5	4926	C	N3-C2-O2	-5.76	117.87	121.90
1	L5	4928	C	C6-N1-C1'	-5.76	113.89	120.80
1	L5	4921	C	N1-C2-O2	5.75	122.35	118.90
1	L5	1417	C	C2-N1-C1'	5.74	125.12	118.80
49	S2	494	C	N1-C2-O2	5.74	122.34	118.90
49	S2	1047	C	C6-N1-C2	-5.74	118.00	120.30
49	S2	1219	C	C2-N1-C1'	5.74	125.11	118.80
72	SG	179	LEU	CA-CB-CG	5.74	128.49	115.30
1	L5	2018	C	C6-N1-C2	-5.73	118.01	120.30
1	L5	4913	G	P-O3'-C3'	5.73	126.58	119.70
1	L5	504	G	N3-C4-C5	-5.69	125.75	128.60
72	SG	189	ARG	CG-CD-NE	5.69	123.75	111.80
1	L5	4398	C	N1-C2-O2	5.69	122.31	118.90
1	L5	4923	C	C6-N1-C2	-5.69	118.03	120.30
49	S2	130	G	N3-C4-N9	5.68	129.41	126.00
1	L5	2094	G	C4-N9-C1'	5.68	133.89	126.50
49	S2	130	G	C4-N9-C1'	5.68	133.89	126.50
1	L5	516	C	N1-C2-O2	5.68	122.31	118.90
1	L5	2532	C	C2-N1-C1'	5.66	125.03	118.80
62	SS	3	LEU	CA-CB-CG	5.66	128.32	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	2494	U	C2-N1-C1'	5.66	124.49	117.70
49	S2	539	C	C2-N1-C1'	5.66	125.03	118.80
49	S2	1271	C	C2-N1-C1'	5.65	125.02	118.80
1	L5	4920	C	N1-C2-O2	5.65	122.29	118.90
49	S2	1417	C	N3-C2-O2	-5.64	117.95	121.90
49	S2	1683	C	N1-C2-O2	5.64	122.29	118.90
83	CA	325	LEU	C-N-CA	5.64	135.80	121.70
1	L5	2416	G	P-O3'-C3'	5.64	126.47	119.70
1	L5	4237	C	C2-N1-C1'	5.63	124.99	118.80
1	L5	2708	U	N3-C2-O2	-5.61	118.27	122.20
49	S2	973	C	N1-C2-O2	5.61	122.27	118.90
1	L5	4147	G	C5-C6-O6	5.61	131.96	128.60
55	SH	32	MET	CB-CG-SD	5.60	129.19	112.40
1	L5	1	C	N1-C2-O2	5.57	122.24	118.90
1	L5	1731	C	C2-N1-C1'	5.56	124.92	118.80
1	L5	1405	C	N3-C2-O2	-5.56	118.01	121.90
1	L5	2528	G	N3-C4-N9	5.55	129.33	126.00
49	S2	126	G	N1-C2-N2	-5.55	111.20	116.20
49	S2	130	G	N3-C4-C5	-5.55	125.82	128.60
51	SB	168	MET	CG-SD-CE	5.55	109.09	100.20
49	S2	583	C	N3-C2-O2	-5.55	118.02	121.90
72	SG	162	LEU	CA-CB-CG	5.55	128.06	115.30
1	L5	2710	C	C6-N1-C2	-5.54	118.08	120.30
1	L5	453	G	N3-C4-N9	5.54	129.32	126.00
25	LW	91	MET	CA-CB-CG	5.54	122.71	113.30
1	L5	4709	U	N3-C2-O2	-5.53	118.33	122.20
60	SQ	52	LEU	CA-CB-CG	5.52	128.00	115.30
1	L5	8	U	C6-N1-C2	-5.52	117.69	121.00
49	S2	593	C	N1-C2-O2	5.52	122.21	118.90
49	S2	1389	C	N1-C2-O2	5.52	122.21	118.90
1	L5	1762	C	N3-C2-O2	-5.51	118.04	121.90
49	S2	1696	C	C2-N1-C1'	5.51	124.86	118.80
49	S2	1865	C	C2-N1-C1'	5.50	124.85	118.80
49	S2	659	G	C4-N9-C1'	5.49	133.64	126.50
49	S2	1111	U	N3-C2-O2	-5.49	118.36	122.20
49	S2	293	C	C6-N1-C1'	-5.48	114.23	120.80
1	L5	2528	G	N3-C4-C5	-5.47	125.86	128.60
1	L5	1472	C	C2-N1-C1'	5.47	124.82	118.80
1	L5	515	C	N3-C2-O2	-5.47	118.07	121.90
49	S2	1309	C	C6-N1-C1'	-5.46	114.24	120.80
56	SI	195	LEU	CB-CG-CD2	-5.46	101.72	111.00
1	L5	3636	C	C6-N1-C2	-5.45	118.12	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4138	C	C6-N1-C2	-5.45	118.12	120.30
1	L5	4709	U	C2-N1-C1'	5.45	124.23	117.70
50	SA	51	LEU	CB-CG-CD2	-5.45	101.74	111.00
21	LS	30	MET	CG-SD-CE	-5.44	91.49	100.20
83	CA	190	LEU	CB-CG-CD2	5.44	120.25	111.00
1	L5	2416	G	OP2-P-O3'	5.44	117.16	105.20
57	SK	79	LEU	CB-CG-CD1	-5.44	101.76	111.00
1	L5	2409	U	C5-C6-N1	5.43	125.42	122.70
26	LX	116	LEU	CA-CB-CG	5.43	127.78	115.30
49	S2	325	C	C2-N1-C1'	5.43	124.77	118.80
71	SC	78	LEU	CA-CB-CG	5.42	127.76	115.30
1	L5	1241	C	C6-N1-C1'	-5.41	114.31	120.80
49	S2	331	C	C2-N1-C1'	5.41	124.75	118.80
49	S2	1434	C	P-O3'-C3'	5.41	126.19	119.70
49	S2	1271	C	N3-C2-O2	-5.40	118.12	121.90
1	L5	4147	G	N1-C6-O6	-5.40	116.66	119.90
1	L5	1633	G	P-O3'-C3'	5.39	126.17	119.70
51	SB	38	MET	CG-SD-CE	-5.39	91.57	100.20
1	L5	1702	C	C2-N1-C1'	5.39	124.73	118.80
1	L5	1977	C	P-O3'-C3'	5.39	126.17	119.70
49	S2	459	C	N1-C2-O2	5.38	122.13	118.90
49	S2	1219	C	N3-C2-O2	-5.38	118.13	121.90
1	L5	1893	C	C2-N1-C1'	5.38	124.72	118.80
2	L7	102	U	N3-C2-O2	-5.38	118.44	122.20
49	S2	112	U	P-O3'-C3'	5.37	126.14	119.70
3	L8	128	C	N3-C4-N4	5.36	121.75	118.00
1	L5	1245	C	C2-N1-C1'	5.35	124.69	118.80
1	L5	4281	A	O4'-C1'-N9	5.35	112.48	108.20
49	S2	850	C	N1-C2-O2	5.35	122.11	118.90
49	S2	1537	A	N7-C8-N9	5.34	116.47	113.80
78	SY	44	LEU	CB-CG-CD1	-5.34	101.93	111.00
1	L5	2410	C	C5-C6-N1	5.33	123.67	121.00
1	L5	1762	C	C6-N1-C2	-5.33	118.17	120.30
49	S2	1309	C	C5-C6-N1	5.33	123.67	121.00
1	L5	2262	G	C8-N9-C1'	-5.32	120.09	127.00
49	S2	291	G	P-O3'-C3'	5.32	126.08	119.70
1	L5	2255	C	N1-C2-O2	5.31	122.08	118.90
1	L5	1816	C	C6-N1-C2	-5.30	118.18	120.30
1	L5	4775	C	C2-N1-C1'	5.30	124.63	118.80
49	S2	579	C	N1-C2-O2	5.30	122.08	118.90
1	L5	5028	G	N3-C4-N9	5.30	129.18	126.00
3	L8	128	C	C6-N1-C2	-5.30	118.18	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	77	U	N1-C2-O2	5.29	126.51	122.80
1	L5	2627	C	N3-C2-O2	-5.29	118.20	121.90
2	L7	28	C	C6-N1-C2	-5.29	118.19	120.30
49	S2	1536	G	C8-N9-C4	-5.29	104.28	106.40
1	L5	3762	U	N1-C2-O2	5.29	126.50	122.80
1	L5	4107	G	N3-C4-N9	5.29	129.17	126.00
49	S2	624	C	N1-C2-O2	5.29	122.07	118.90
55	SH	40	LEU	CA-CB-CG	5.28	127.45	115.30
1	L5	406	C	P-O3'-C3'	5.28	126.03	119.70
1	L5	2528	G	C8-N9-C1'	-5.27	120.15	127.00
79	SZ	47	LEU	CA-CB-CG	5.27	127.41	115.30
1	L5	1252	C	C6-N1-C2	-5.26	118.19	120.30
1	L5	1241	C	C6-N1-C2	-5.26	118.19	120.30
1	L5	1412	G	N7-C8-N9	5.26	115.73	113.10
49	S2	201	C	N1-C2-O2	5.26	122.06	118.90
49	S2	530	U	N1-C2-O2	5.26	126.48	122.80
49	S2	1520	G	N3-C4-C5	-5.26	125.97	128.60
1	L5	1726	U	N1-C2-O2	5.25	126.48	122.80
1	L5	2786	C	P-O3'-C3'	5.25	126.00	119.70
1	L5	1577	G	N3-C2-N2	-5.23	116.24	119.90
49	S2	530	U	C2-N1-C1'	5.23	123.97	117.70
1	L5	3778	U	N1-C2-O2	5.22	126.46	122.80
1	L5	3673	C	P-O3'-C3'	5.22	125.96	119.70
1	L5	4572	U	N1-C2-O2	5.21	126.45	122.80
49	S2	1660	C	C2-N1-C1'	5.21	124.53	118.80
49	S2	1261	C	C2-N1-C1'	5.21	124.53	118.80
1	L5	3771	C	C2-N1-C1'	5.21	124.53	118.80
1	L5	2094	G	N3-C4-N9	5.20	129.12	126.00
49	S2	1263	U	C2-N3-C4	-5.20	123.88	127.00
49	S2	331	C	N3-C4-C5	-5.20	119.82	121.90
49	S2	798	A	C8-N9-C4	5.20	107.88	105.80
1	L5	3909	C	C2-N1-C1'	5.19	124.51	118.80
1	L5	1182	C	N3-C2-O2	-5.19	118.27	121.90
1	L5	2018	C	C2-N1-C1'	5.18	124.50	118.80
1	L5	1735	U	N1-C2-O2	5.17	126.42	122.80
49	S2	530	U	N3-C2-O2	-5.17	118.58	122.20
1	L5	3882	C	C2-N1-C1'	5.15	124.47	118.80
1	L5	1853	G	C4-N9-C1'	5.15	133.20	126.50
1	L5	294	G	C4-N9-C1'	5.14	133.18	126.50
1	L5	2820	C	N3-C2-O2	-5.13	118.31	121.90
1	L5	2627	C	C5-C6-N1	5.13	123.56	121.00
1	L5	515	C	C6-N1-C1'	-5.12	114.66	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	1809	C	C2-N1-C1'	5.12	124.43	118.80
49	S2	1315	U	N3-C2-O2	-5.11	118.62	122.20
1	L5	2410	C	C6-N1-C2	-5.11	118.26	120.30
6	LC	319	LEU	CB-CG-CD1	-5.10	102.33	111.00
1	L5	4148	C	N1-C2-O2	5.10	121.96	118.90
1	L5	4682	U	N3-C2-O2	-5.10	118.63	122.20
1	L5	4928	C	O4'-C1'-N1	5.10	112.28	108.20
1	L5	4360	U	N3-C2-O2	-5.09	118.63	122.20
49	S2	118	C	N1-C2-O2	5.09	121.96	118.90
1	L5	220	C	C6-N1-C2	-5.09	118.26	120.30
49	S2	180	G	N3-C4-C5	-5.09	126.05	128.60
1	L5	4901	G	N1-C6-O6	-5.08	116.85	119.90
1	L5	1340	C	C5-C6-N1	5.08	123.54	121.00
1	L5	4101	C	N3-C4-C5	5.08	123.93	121.90
1	L5	485	C	N3-C2-O2	-5.08	118.34	121.90
49	S2	1520	G	C6-C5-N7	-5.08	127.35	130.40
1	L5	4107	G	C8-N9-C1'	-5.07	120.41	127.00
49	S2	965	U	N3-C2-O2	-5.07	118.65	122.20
1	L5	673	C	C5-C6-N1	5.07	123.53	121.00
49	S2	1218	C	C5-C6-N1	5.06	123.53	121.00
1	L5	1193	C	C6-N1-C2	-5.06	118.28	120.30
49	S2	126	G	N3-C2-N2	5.06	123.44	119.90
49	S2	325	C	N1-C2-O2	5.06	121.94	118.90
1	L5	4924	C	N1-C2-O2	5.06	121.93	118.90
1	L5	4747	C	C2-N1-C1'	5.05	124.36	118.80
55	SH	35	ASP	CA-C-O	5.05	130.71	120.10
49	S2	1746	U	N3-C2-O2	-5.05	118.67	122.20
1	L5	2627	C	C6-N1-C1'	-5.05	114.74	120.80
23	LU	43	LEU	CA-CB-CG	5.04	126.90	115.30
49	S2	1292	C	C2-N1-C1'	5.04	124.35	118.80
49	S2	427	U	N3-C2-O2	-5.04	118.67	122.20
49	S2	1230	C	C6-N1-C2	-5.04	118.28	120.30
1	L5	1273	G	C4-N9-C1'	5.04	133.05	126.50
1	L5	3693	U	N3-C2-O2	-5.04	118.67	122.20
1	L5	4241	C	C2-N1-C1'	5.03	124.33	118.80
1	L5	1259	G	N1-C6-O6	-5.02	116.89	119.90
1	L5	4773	C	C2-N1-C1'	5.01	124.32	118.80
49	S2	4	C	C2-N1-C1'	5.01	124.31	118.80
1	L5	1273	G	C8-N9-C1'	-5.01	120.48	127.00
1	L5	2900	U	C2-N1-C1'	5.01	123.71	117.70

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	LA	110	GLY	Peptide
13	LJ	18	ARG	Peptide
17	LO	30	GLY	Peptide
25	LW	19	ARG	Sidechain
30	Lb	117	ARG	Sidechain
35	Lg	29	ARG	Sidechain
37	Li	56	ARG	Sidechain
50	SA	205	ARG	Sidechain
53	SE	66	MET	Peptide
54	SF	78	MET	Peptide
72	SG	189	ARG	Sidechain
62	SS	23	ARG	Sidechain
62	SS	38	ARG	Sidechain

5.2 Too-close contacts [i](#)

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

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	LA	246/257 (96%)	227 (92%)	19 (8%)	0	100	100
5	LB	400/403 (99%)	379 (95%)	20 (5%)	1 (0%)	41	76
6	LC	366/427 (86%)	340 (93%)	26 (7%)	0	100	100
7	LD	291/297 (98%)	277 (95%)	14 (5%)	0	100	100
8	LE	215/288 (75%)	199 (93%)	16 (7%)	0	100	100
9	LF	223/248 (90%)	215 (96%)	8 (4%)	0	100	100
10	LG	239/266 (90%)	225 (94%)	14 (6%)	0	100	100
11	LH	188/192 (98%)	172 (92%)	16 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
12	LI	198/214 (92%)	187 (94%)	10 (5%)	1 (0%)	29	68
13	LJ	174/178 (98%)	161 (92%)	13 (8%)	0	100	100
14	LL	208/211 (99%)	196 (94%)	12 (6%)	0	100	100
15	LM	137/215 (64%)	128 (93%)	9 (7%)	0	100	100
16	LN	201/204 (98%)	190 (94%)	10 (5%)	1 (0%)	29	68
17	LO	199/203 (98%)	194 (98%)	5 (2%)	0	100	100
18	LP	151/184 (82%)	142 (94%)	9 (6%)	0	100	100
19	LQ	185/188 (98%)	176 (95%)	9 (5%)	0	100	100
20	LR	185/196 (94%)	181 (98%)	4 (2%)	0	100	100
21	LS	173/176 (98%)	161 (93%)	12 (7%)	0	100	100
22	LT	157/160 (98%)	151 (96%)	6 (4%)	0	100	100
23	LU	98/128 (77%)	93 (95%)	5 (5%)	0	100	100
24	LV	129/140 (92%)	122 (95%)	7 (5%)	0	100	100
25	LW	122/157 (78%)	113 (93%)	9 (7%)	0	100	100
26	LX	118/156 (76%)	116 (98%)	2 (2%)	0	100	100
27	LY	131/145 (90%)	126 (96%)	5 (4%)	0	100	100
28	LZ	133/136 (98%)	127 (96%)	6 (4%)	0	100	100
29	La	145/148 (98%)	137 (94%)	8 (6%)	0	100	100
30	Lb	105/159 (66%)	98 (93%)	7 (7%)	0	100	100
31	Lc	96/115 (84%)	92 (96%)	4 (4%)	0	100	100
32	Ld	105/125 (84%)	97 (92%)	8 (8%)	0	100	100
33	Le	126/135 (93%)	122 (97%)	4 (3%)	0	100	100
34	Lf	107/110 (97%)	100 (94%)	7 (6%)	0	100	100
35	Lg	112/117 (96%)	111 (99%)	1 (1%)	0	100	100
36	Lh	120/123 (98%)	119 (99%)	1 (1%)	0	100	100
37	Li	100/105 (95%)	100 (100%)	0	0	100	100
38	Lj	84/97 (87%)	76 (90%)	8 (10%)	0	100	100
39	Lk	67/70 (96%)	63 (94%)	4 (6%)	0	100	100
40	Ll	48/51 (94%)	47 (98%)	1 (2%)	0	100	100
41	Lm	49/128 (38%)	49 (100%)	0	0	100	100
42	Ln	22/25 (88%)	22 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
43	Lo	103/106 (97%)	97 (94%)	5 (5%)	1 (1%)	15	53
44	Lp	89/92 (97%)	83 (93%)	6 (7%)	0	100	100
45	Lr	123/137 (90%)	117 (95%)	6 (5%)	0	100	100
46	Ls	194/317 (61%)	174 (90%)	20 (10%)	0	100	100
47	Lt	137/165 (83%)	107 (78%)	29 (21%)	1 (1%)	22	60
48	Lz	203/217 (94%)	180 (89%)	23 (11%)	0	100	100
50	SA	219/295 (74%)	208 (95%)	11 (5%)	0	100	100
51	SB	212/264 (80%)	202 (95%)	10 (5%)	0	100	100
52	SD	225/243 (93%)	207 (92%)	18 (8%)	0	100	100
53	SE	260/263 (99%)	242 (93%)	17 (6%)	1 (0%)	34	72
54	SF	180/204 (88%)	172 (96%)	7 (4%)	1 (1%)	25	64
55	SH	182/194 (94%)	173 (95%)	9 (5%)	0	100	100
56	SI	204/208 (98%)	197 (97%)	7 (3%)	0	100	100
57	SK	93/165 (56%)	85 (91%)	8 (9%)	0	100	100
58	SL	140/158 (89%)	134 (96%)	6 (4%)	0	100	100
59	SP	127/145 (88%)	119 (94%)	8 (6%)	0	100	100
60	SQ	142/146 (97%)	132 (93%)	10 (7%)	0	100	100
61	SR	133/135 (98%)	126 (95%)	7 (5%)	0	100	100
62	SS	142/152 (93%)	133 (94%)	9 (6%)	0	100	100
63	ST	141/145 (97%)	133 (94%)	7 (5%)	1 (1%)	22	60
64	SU	102/119 (86%)	93 (91%)	9 (9%)	0	100	100
65	SV	81/83 (98%)	79 (98%)	2 (2%)	0	100	100
66	SX	139/143 (97%)	129 (93%)	9 (6%)	1 (1%)	22	60
67	Sa	100/115 (87%)	91 (91%)	8 (8%)	1 (1%)	15	53
68	Sc	62/69 (90%)	59 (95%)	3 (5%)	0	100	100
69	Sd	53/56 (95%)	49 (92%)	3 (6%)	1 (2%)	8	36
70	Sg	311/317 (98%)	275 (88%)	36 (12%)	0	100	100
71	SC	220/293 (75%)	207 (94%)	13 (6%)	0	100	100
72	SG	235/249 (94%)	224 (95%)	11 (5%)	0	100	100
73	SJ	183/194 (94%)	173 (94%)	9 (5%)	1 (0%)	29	68
74	SM	120/132 (91%)	113 (94%)	7 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
75	SN	148/151 (98%)	143 (97%)	5 (3%)	0	100	100
76	SO	133/151 (88%)	120 (90%)	12 (9%)	1 (1%)	19	57
77	SW	127/130 (98%)	120 (94%)	7 (6%)	0	100	100
78	SY	124/133 (93%)	118 (95%)	6 (5%)	0	100	100
79	SZ	73/125 (58%)	58 (80%)	15 (20%)	0	100	100
80	Sb	81/84 (96%)	73 (90%)	8 (10%)	0	100	100
81	Se	56/59 (95%)	50 (89%)	6 (11%)	0	100	100
82	Sf	65/156 (42%)	50 (77%)	15 (23%)	0	100	100
83	CA	350/394 (89%)	332 (95%)	18 (5%)	0	100	100
84	CB	14/408 (3%)	14 (100%)	0	0	100	100
86	CE	71/223 (32%)	70 (99%)	1 (1%)	0	100	100
All	All	12250/14412 (85%)	11492 (94%)	745 (6%)	13 (0%)	54	85

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
12	LI	213	HIS
43	Lo	56	PHE
5	LB	302	ASN
16	LN	124	ASP
53	SE	165	GLU
54	SF	128	ILE
67	Sa	46	GLU
73	SJ	122	SER
63	ST	45	LEU
66	SX	124	LYS
76	SO	140	THR
69	Sd	14	PHE
47	Lt	9	GLU

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	LA	190/199 (96%)	190 (100%)	0	100	100
5	LB	348/349 (100%)	347 (100%)	1 (0%)	92	97
6	LC	306/348 (88%)	305 (100%)	1 (0%)	92	97
7	LD	246/250 (98%)	246 (100%)	0	100	100
8	LE	195/252 (77%)	193 (99%)	2 (1%)	76	91
9	LF	194/215 (90%)	194 (100%)	0	100	100
10	LG	203/223 (91%)	203 (100%)	0	100	100
11	LH	169/171 (99%)	169 (100%)	0	100	100
12	LI	172/181 (95%)	172 (100%)	0	100	100
13	LJ	148/149 (99%)	147 (99%)	1 (1%)	84	94
14	LL	176/177 (99%)	176 (100%)	0	100	100
15	LM	118/161 (73%)	118 (100%)	0	100	100
16	LN	171/172 (99%)	171 (100%)	0	100	100
17	LO	173/174 (99%)	173 (100%)	0	100	100
18	LP	134/163 (82%)	132 (98%)	2 (2%)	65	87
19	LQ	164/165 (99%)	164 (100%)	0	100	100
20	LR	166/175 (95%)	166 (100%)	0	100	100
21	LS	156/157 (99%)	156 (100%)	0	100	100
22	LT	139/140 (99%)	139 (100%)	0	100	100
23	LU	90/115 (78%)	89 (99%)	1 (1%)	73	90
24	LV	101/107 (94%)	101 (100%)	0	100	100
25	LW	103/126 (82%)	103 (100%)	0	100	100
26	LX	108/133 (81%)	108 (100%)	0	100	100
27	LY	123/135 (91%)	123 (100%)	0	100	100
28	LZ	117/118 (99%)	117 (100%)	0	100	100
29	La	120/121 (99%)	120 (100%)	0	100	100
30	Lb	89/126 (71%)	89 (100%)	0	100	100
31	Lc	83/97 (86%)	83 (100%)	0	100	100
32	Ld	98/110 (89%)	98 (100%)	0	100	100
33	Le	114/121 (94%)	114 (100%)	0	100	100
34	Lf	88/89 (99%)	88 (100%)	0	100	100
35	Lg	98/100 (98%)	98 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
36	Lh	109/110 (99%)	109 (100%)	0	100	100
37	Li	86/89 (97%)	85 (99%)	1 (1%)	71	90
38	Lj	73/80 (91%)	73 (100%)	0	100	100
39	Lk	64/65 (98%)	64 (100%)	0	100	100
40	Ll	47/48 (98%)	46 (98%)	1 (2%)	53	82
41	Lm	47/115 (41%)	47 (100%)	0	100	100
42	Ln	23/24 (96%)	23 (100%)	0	100	100
43	Lo	93/94 (99%)	92 (99%)	1 (1%)	73	90
44	Lp	74/75 (99%)	74 (100%)	0	100	100
45	Lr	109/121 (90%)	109 (100%)	0	100	100
46	Ls	162/258 (63%)	162 (100%)	0	100	100
47	Lt	112/137 (82%)	111 (99%)	1 (1%)	78	92
50	SA	183/243 (75%)	183 (100%)	0	100	100
51	SB	195/231 (84%)	195 (100%)	0	100	100
52	SD	190/202 (94%)	189 (100%)	1 (0%)	88	96
53	SE	224/225 (100%)	224 (100%)	0	100	100
54	SF	156/170 (92%)	156 (100%)	0	100	100
55	SH	166/174 (95%)	166 (100%)	0	100	100
56	SI	178/180 (99%)	178 (100%)	0	100	100
57	SK	86/136 (63%)	86 (100%)	0	100	100
58	SL	130/142 (92%)	129 (99%)	1 (1%)	81	93
59	SP	115/130 (88%)	113 (98%)	2 (2%)	60	85
60	SQ	119/121 (98%)	119 (100%)	0	100	100
61	SR	122/122 (100%)	121 (99%)	1 (1%)	81	93
62	SS	125/132 (95%)	125 (100%)	0	100	100
63	ST	113/115 (98%)	113 (100%)	0	100	100
64	SU	94/107 (88%)	94 (100%)	0	100	100
65	SV	67/67 (100%)	67 (100%)	0	100	100
66	SX	113/115 (98%)	113 (100%)	0	100	100
67	Sa	89/98 (91%)	89 (100%)	0	100	100
68	Sc	57/62 (92%)	57 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
69	Sd	48/49 (98%)	48 (100%)	0	100	100
70	Sg	272/275 (99%)	271 (100%)	1 (0%)	91	97
71	SC	188/225 (84%)	187 (100%)	1 (0%)	88	96
72	SG	207/218 (95%)	203 (98%)	4 (2%)	57	84
73	SJ	161/168 (96%)	161 (100%)	0	100	100
75	SN	130/131 (99%)	130 (100%)	0	100	100
76	SO	105/119 (88%)	105 (100%)	0	100	100
77	SW	112/113 (99%)	111 (99%)	1 (1%)	78	92
78	SY	109/115 (95%)	109 (100%)	0	100	100
79	SZ	66/103 (64%)	66 (100%)	0	100	100
80	Sb	75/76 (99%)	75 (100%)	0	100	100
81	Se	47/48 (98%)	46 (98%)	1 (2%)	53	82
82	Sf	60/140 (43%)	56 (93%)	4 (7%)	16	49
83	CA	303/336 (90%)	302 (100%)	1 (0%)	92	97
84	CB	14/328 (4%)	14 (100%)	0	100	100
86	CE	62/190 (33%)	62 (100%)	0	100	100
All	All	10380/11941 (87%)	10350 (100%)	30 (0%)	92	97

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	LB	361	GLU
6	LC	188	ARG
8	LE	56	ARG
8	LE	100	LYS
13	LJ	58	ARG
18	LP	10	ASN
18	LP	69	ARG
23	LU	67	LYS
37	Li	56	ARG
40	Ll	36	ARG
43	Lo	99	ARG
47	Lt	123	ARG
52	SD	76	ARG
58	SL	69	ARG
59	SP	14	LYS

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Mol	Chain	Res	Type
59	SP	127	LYS
61	SR	72	LYS
70	Sg	118	ARG
71	SC	167	ARG
72	SG	142	ARG
72	SG	189	ARG
72	SG	200	LYS
72	SG	211	LYS
77	SW	15	ASN
81	Se	24	LYS
82	Sf	89	LYS
82	Sf	104	LYS
82	Sf	118	ARG
82	Sf	138	ARG
83	CA	71	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
5	LB	289	GLN
52	SD	174	HIS
59	SP	24	GLN
61	SR	48	ASN
63	ST	137	GLN
66	SX	63	ASN
68	Sc	7	GLN
71	SC	235	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	L5	3704/5070 (73%)	814 (21%)	20 (0%)
2	L7	119/121 (98%)	9 (7%)	0
3	L8	155/157 (98%)	26 (16%)	0
49	S2	1717/1869 (91%)	452 (26%)	5 (0%)
85	CC	67/75 (89%)	17 (25%)	1 (1%)
All	All	5762/7292 (79%)	1318 (22%)	26 (0%)

All (1318) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	L5	17	A
1	L5	25	A
1	L5	30	C
1	L5	39	A
1	L5	42	A
1	L5	48	G
1	L5	56	A
1	L5	59	A
1	L5	64	A
1	L5	65	A
1	L5	66	A
1	L5	69	A
1	L5	73	A
1	L5	91	G
1	L5	98	A
1	L5	104	G
1	L5	108	A
1	L5	109	G
1	L5	110	C
1	L5	112	C
1	L5	119	G
1	L5	120	A
1	L5	127	G
1	L5	128	C
1	L5	132	G
1	L5	133	C
1	L5	134	G
1	L5	135	G
1	L5	136	C
1	L5	137	G
1	L5	141	C
1	L5	152	U
1	L5	159	C
1	L5	171	U
1	L5	172	C
1	L5	173	C
1	L5	177	G
1	L5	181	C
1	L5	183	C
1	L5	184	U
1	L5	185	C
1	L5	186	G
1	L5	187	U

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Mol	Chain	Res	Type
1	L5	188	G
1	L5	189	G
1	L5	200	U
1	L5	210	C
1	L5	217	C
1	L5	218	A
1	L5	219	G
1	L5	220	C
1	L5	233	U
1	L5	234	G
1	L5	236	G
1	L5	261	G
1	L5	264	C
1	L5	265	C
1	L5	266	C
1	L5	280	G
1	L5	297	U
1	L5	306	A
1	L5	315	G
1	L5	316	U
1	L5	340	C
1	L5	360	A
1	L5	373	G
1	L5	387	G
1	L5	407	A
1	L5	409	G
1	L5	410	A
1	L5	411	G
1	L5	412	G
1	L5	413	G
1	L5	431	G
1	L5	432	U
1	L5	438	G
1	L5	440	U
1	L5	449	C
1	L5	450	G
1	L5	452	A
1	L5	453	G
1	L5	454	U
1	L5	465	G
1	L5	467	U
1	L5	468	U

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Mol	Chain	Res	Type
1	L5	484	U
1	L5	485	C
1	L5	486	C
1	L5	489	C
1	L5	493	G
1	L5	494	U
1	L5	495	C
1	L5	497	G
1	L5	498	C
1	L5	500	G
1	L5	502	C
1	L5	503	C
1	L5	504	G
1	L5	505	G
1	L5	506	C
1	L5	509	A
1	L5	510	U
1	L5	512	U
1	L5	513	U
1	L5	514	U
1	L5	515	C
1	L5	516	C
1	L5	517	C
1	L5	518	G
1	L5	643	C
1	L5	646	G
1	L5	657	C
1	L5	659	G
1	L5	665	C
1	L5	666	G
1	L5	667	A
1	L5	668	C
1	L5	669	C
1	L5	670	G
1	L5	672	C
1	L5	673	C
1	L5	684	G
1	L5	685	C
1	L5	686	A
1	L5	687	U
1	L5	696	C
1	L5	703	G

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Mol	Chain	Res	Type
1	L5	704	C
1	L5	708	G
1	L5	731	G
1	L5	738	C
1	L5	739	G
1	L5	740	G
1	L5	742	G
1	L5	757	G
1	L5	758	G
1	L5	759	G
1	L5	760	G
1	L5	904	C
1	L5	906	C
1	L5	913	U
1	L5	914	U
1	L5	915	A
1	L5	917	A
1	L5	918	G
1	L5	923	C
1	L5	924	C
1	L5	926	G
1	L5	932	A
1	L5	933	G
1	L5	935	A
1	L5	936	C
1	L5	941	C
1	L5	943	A
1	L5	945	U
1	L5	956	A
1	L5	959	G
1	L5	960	A
1	L5	961	G
1	L5	962	C
1	L5	965	G
1	L5	966	A
1	L5	967	C
1	L5	969	C
1	L5	970	G
1	L5	972	C
1	L5	982	U
1	L5	989	U
1	L5	990	C

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Mol	Chain	Res	Type
1	L5	992	C
1	L5	993	G
1	L5	995	C
1	L5	996	G
1	L5	1048	G
1	L5	1049	C
1	L5	1050	C
1	L5	1051	G
1	L5	1070	G
1	L5	1071	C
1	L5	1072	C
1	L5	1075	G
1	L5	1095	A
1	L5	1168	G
1	L5	1171	G
1	L5	1172	C
1	L5	1173	G
1	L5	1178	G
1	L5	1179	U
1	L5	1180	C
1	L5	1181	C
1	L5	1182	C
1	L5	1183	C
1	L5	1193	C
1	L5	1202	C
1	L5	1203	G
1	L5	1204	C
1	L5	1210	C
1	L5	1211	G
1	L5	1214	C
1	L5	1215	C
1	L5	1216	C
1	L5	1217	G
1	L5	1219	G
1	L5	1222	A
1	L5	1235	G
1	L5	1246	G
1	L5	1247	U
1	L5	1253	G
1	L5	1254	A
1	L5	1257	A
1	L5	1258	G

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Mol	Chain	Res	Type
1	L5	1266	G
1	L5	1267	C
1	L5	1271	G
1	L5	1272	C
1	L5	1273	G
1	L5	1274	A
1	L5	1275	G
1	L5	1280	C
1	L5	1284	G
1	L5	1285	U
1	L5	1287	G
1	L5	1294	A
1	L5	1295	C
1	L5	1296	G
1	L5	1301	C
1	L5	1302	U
1	L5	1324	A
1	L5	1326	A
1	L5	1337	A
1	L5	1344	C
1	L5	1354	A
1	L5	1358	G
1	L5	1359	G
1	L5	1365	C
1	L5	1366	G
1	L5	1387	A
1	L5	1393	G
1	L5	1394	G
1	L5	1397	A
1	L5	1407	C
1	L5	1408	G
1	L5	1409	C
1	L5	1410	U
1	L5	1411	C
1	L5	1414	C
1	L5	1416	G
1	L5	1420	A
1	L5	1435	G
1	L5	1438	U
1	L5	1444	G
1	L5	1447	C
1	L5	1457	G

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Mol	Chain	Res	Type
1	L5	1482	G
1	L5	1483	C
1	L5	1497	A
1	L5	1498	G
1	L5	1502	G
1	L5	1504	G
1	L5	1517	G
1	L5	1518	A
1	L5	1525	A
1	L5	1534	A
1	L5	1535	C
1	L5	1547	A
1	L5	1562	G
1	L5	1566	C
1	L5	1574	G
1	L5	1578	U
1	L5	1591	U
1	L5	1596	U
1	L5	1613	A
1	L5	1614	C
1	L5	1624	G
1	L5	1625	G
1	L5	1631	A
1	L5	1633	G
1	L5	1634	A
1	L5	1637	A
1	L5	1638	A
1	L5	1640	C
1	L5	1641	G
1	L5	1642	A
1	L5	1654	G
1	L5	1661	C
1	L5	1670	G
1	L5	1676	C
1	L5	1677	U
1	L5	1678	C
1	L5	1680	G
1	L5	1691	G
1	L5	1694	C
1	L5	1697	G
1	L5	1699	A
1	L5	1700	G

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Mol	Chain	Res	Type
1	L5	1703	C
1	L5	1704	C
1	L5	1705	G
1	L5	1708	G
1	L5	1718	C
1	L5	1719	A
1	L5	1726	U
1	L5	1731	C
1	L5	1734	G
1	L5	1741	G
1	L5	1742	A
1	L5	1753	G
1	L5	1755	C
1	L5	1757	U
1	L5	1758	G
1	L5	1760	G
1	L5	1761	G
1	L5	1762	C
1	L5	1763	C
1	L5	1764	G
1	L5	1765	A
1	L5	1766	A
1	L5	1768	C
1	L5	1770	A
1	L5	1771	U
1	L5	1772	C
1	L5	1775	A
1	L5	1785	C
1	L5	1787	A
1	L5	1804	A
1	L5	1810	G
1	L5	1815	G
1	L5	1820	C
1	L5	1821	G
1	L5	1822	U
1	L5	1834	U
1	L5	1836	G
1	L5	1837	A
1	L5	1842	G
1	L5	1855	G
1	L5	1869	G
1	L5	1882	U

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Mol	Chain	Res	Type
1	L5	1892	A
1	L5	1897	A
1	L5	1898	C
1	L5	1917	A
1	L5	1918	U
1	L5	1919	G
1	L5	1920	C
1	L5	1921	C
1	L5	1922	G
1	L5	1925	G
1	L5	1930	U
1	L5	1931	C
1	L5	1932	A
1	L5	1936	C
1	L5	1940	G
1	L5	1946	G
1	L5	1948	G
1	L5	1949	U
1	L5	1951	G
1	L5	1959	U
1	L5	1960	A
1	L5	1961	G
1	L5	1962	A
1	L5	1972	G
1	L5	1974	U
1	L5	1975	G
1	L5	1978	C
1	L5	1980	U
1	L5	1983	A
1	L5	1984	A
1	L5	1985	G
1	L5	1986	U
1	L5	1987	C
1	L5	1988	G
1	L5	1990	A
1	L5	1992	U
1	L5	1993	C
1	L5	1997	U
1	L5	2001	G
1	L5	2002	A
1	L5	2004	U
1	L5	2014	C

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Mol	Chain	Res	Type
1	L5	2017	A
1	L5	2018	C
1	L5	2024	G
1	L5	2026	A
1	L5	2034	G
1	L5	2046	G
1	L5	2048	U
1	L5	2055	G
1	L5	2056	G
1	L5	2069	A
1	L5	2084	C
1	L5	2085	G
1	L5	2089	G
1	L5	2092	G
1	L5	2093	A
1	L5	2095	A
1	L5	2096	G
1	L5	2097	U
1	L5	2098	G
1	L5	2102	G
1	L5	2106	G
1	L5	2107	C
1	L5	2108	G
1	L5	2110	C
1	L5	2111	G
1	L5	2112	G
1	L5	2250	C
1	L5	2252	G
1	L5	2253	A
1	L5	2256	C
1	L5	2289	C
1	L5	2300	A
1	L5	2301	G
1	L5	2306	G
1	L5	2313	A
1	L5	2314	G
1	L5	2331	G
1	L5	2332	A
1	L5	2333	G
1	L5	2345	G
1	L5	2348	G
1	L5	2351	C

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Mol	Chain	Res	Type
1	L5	2360	A
1	L5	2369	U
1	L5	2395	A
1	L5	2397	G
1	L5	2408	U
1	L5	2417	A
1	L5	2418	A
1	L5	2421	G
1	L5	2425	U
1	L5	2450	G
1	L5	2464	C
1	L5	2465	C
1	L5	2467	U
1	L5	2474	G
1	L5	2475	G
1	L5	2478	C
1	L5	2479	G
1	L5	2484	A
1	L5	2485	U
1	L5	2487	G
1	L5	2488	C
1	L5	2489	C
1	L5	2490	U
1	L5	2491	C
1	L5	2494	U
1	L5	2495	U
1	L5	2503	G
1	L5	2504	C
1	L5	2505	C
1	L5	2506	G
1	L5	2513	A
1	L5	2519	U
1	L5	2529	A
1	L5	2537	A
1	L5	2544	G
1	L5	2546	G
1	L5	2547	G
1	L5	2554	U
1	L5	2555	G
1	L5	2559	G
1	L5	2560	C
1	L5	2565	A

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Mol	Chain	Res	Type
1	L5	2573	A
1	L5	2583	C
1	L5	2587	A
1	L5	2589	C
1	L5	2601	A
1	L5	2611	A
1	L5	2627	C
1	L5	2653	C
1	L5	2662	G
1	L5	2669	C
1	L5	2670	C
1	L5	2676	A
1	L5	2687	U
1	L5	2694	G
1	L5	2695	A
1	L5	2696	A
1	L5	2703	G
1	L5	2707	U
1	L5	2708	U
1	L5	2709	C
1	L5	2710	C
1	L5	2711	G
1	L5	2721	G
1	L5	2724	G
1	L5	2726	G
1	L5	2738	C
1	L5	2739	C
1	L5	2742	G
1	L5	2743	A
1	L5	2756	G
1	L5	2761	U
1	L5	2763	U
1	L5	2764	A
1	L5	2770	C
1	L5	2787	A
1	L5	2788	U
1	L5	2790	U
1	L5	2814	C
1	L5	2815	A
1	L5	2825	A
1	L5	2826	U
1	L5	2827	G

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Mol	Chain	Res	Type
1	L5	2838	G
1	L5	2845	A
1	L5	2848	G
1	L5	2855	G
1	L5	2877	G
1	L5	2892	C
1	L5	2900	U
1	L5	2902	G
1	L5	2903	G
1	L5	2904	U
1	L5	2905	C
1	L5	2906	G
1	L5	2908	U
1	L5	3585	G
1	L5	3587	C
1	L5	3588	C
1	L5	3590	G
1	L5	3591	C
1	L5	3594	C
1	L5	3595	U
1	L5	3596	A
1	L5	3597	G
1	L5	3604	A
1	L5	3606	U
1	L5	3615	G
1	L5	3616	U
1	L5	3618	C
1	L5	3626	G
1	L5	3630	A
1	L5	3635	A
1	L5	3644	U
1	L5	3646	A
1	L5	3648	A
1	L5	3662	A
1	L5	3673	C
1	L5	3674	G
1	L5	3710	G
1	L5	3714	G
1	L5	3727	A
1	L5	3729	U
1	L5	3735	G
1	L5	3736	A

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Mol	Chain	Res	Type
1	L5	3750	G
1	L5	3753	G
1	L5	3757	G
1	L5	3758	U
1	L5	3759	A
1	L5	3760	A
1	L5	3761	C
1	L5	3771	C
1	L5	3776	G
1	L5	3777	G
1	L5	3784	A
1	L5	3786	U
1	L5	3802	U
1	L5	3811	G
1	L5	3812	C
1	L5	3814	U
1	L5	3817	A
1	L5	3818	U
1	L5	3819	G
1	L5	3823	G
1	L5	3838	U
1	L5	3839	G
1	L5	3840	U
1	L5	3867	A
1	L5	3877	A
1	L5	3878	C
1	L5	3879	G
1	L5	3885	G
1	L5	3887	C
1	L5	3890	A
1	L5	3892	U
1	L5	3897	G
1	L5	3901	A
1	L5	3906	A
1	L5	3907	G
1	L5	3908	A
1	L5	3915	U
1	L5	3926	C
1	L5	3938	G
1	L5	3939	G
1	L5	3942	A
1	L5	3943	A

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Mol	Chain	Res	Type
1	L5	3947	A
1	L5	3950	U
1	L5	3951	G
1	L5	3955	G
1	L5	3956	G
1	L5	3958	G
1	L5	3959	U
1	L5	3960	A
1	L5	3961	G
1	L5	3962	A
1	L5	3963	A
1	L5	3964	U
1	L5	3965	A
1	L5	3966	A
1	L5	3969	G
1	L5	3970	G
1	L5	3971	G
1	L5	3973	G
1	L5	3974	G
1	L5	3975	C
1	L5	3977	C
1	L5	4034	G
1	L5	4035	G
1	L5	4036	G
1	L5	4038	C
1	L5	4039	G
1	L5	4041	C
1	L5	4042	G
1	L5	4043	G
1	L5	4044	U
1	L5	4045	G
1	L5	4046	A
1	L5	4048	A
1	L5	4049	U
1	L5	4051	C
1	L5	4052	C
1	L5	4053	A
1	L5	4055	U
1	L5	4064	C
1	L5	4065	G
1	L5	4076	G
1	L5	4084	G

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Mol	Chain	Res	Type
1	L5	4086	G
1	L5	4097	G
1	L5	4099	G
1	L5	4101	C
1	L5	4102	C
1	L5	4104	G
1	L5	4106	G
1	L5	4107	G
1	L5	4108	G
1	L5	4111	U
1	L5	4112	C
1	L5	4113	U
1	L5	4114	C
1	L5	4115	G
1	L5	4116	C
1	L5	4117	U
1	L5	4119	C
1	L5	4120	U
1	L5	4122	G
1	L5	4127	A
1	L5	4133	C
1	L5	4140	C
1	L5	4141	G
1	L5	4142	C
1	L5	4143	G
1	L5	4144	C
1	L5	4146	G
1	L5	4149	C
1	L5	4150	G
1	L5	4162	C
1	L5	4163	U
1	L5	4168	G
1	L5	4170	A
1	L5	4183	G
1	L5	4191	G
1	L5	4196	G
1	L5	4197	G
1	L5	4203	A
1	L5	4222	G
1	L5	4228	G
1	L5	4229	U
1	L5	4233	A

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Mol	Chain	Res	Type
1	L5	4242	U
1	L5	4251	A
1	L5	4254	G
1	L5	4257	A
1	L5	4265	U
1	L5	4268	A
1	L5	4273	A
1	L5	4291	G
1	L5	4304	A
1	L5	4305	G
1	L5	4314	C
1	L5	4329	G
1	L5	4330	G
1	L5	4332	C
1	L5	4349	C
1	L5	4354	U
1	L5	4373	G
1	L5	4376	A
1	L5	4377	G
1	L5	4378	A
1	L5	4379	A
1	L5	4380	A
1	L5	4387	C
1	L5	4391	G
1	L5	4394	A
1	L5	4422	A
1	L5	4426	C
1	L5	4448	G
1	L5	4449	A
1	L5	4453	C
1	L5	4464	A
1	L5	4475	G
1	L5	4488	A
1	L5	4500	U
1	L5	4512	U
1	L5	4513	A
1	L5	4519	C
1	L5	4524	G
1	L5	4545	G
1	L5	4548	A
1	L5	4549	G
1	L5	4560	C

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Mol	Chain	Res	Type
1	L5	4567	G
1	L5	4573	G
1	L5	4575	G
1	L5	4584	A
1	L5	4590	A
1	L5	4600	G
1	L5	4617	G
1	L5	4630	G
1	L5	4636	U
1	L5	4637	G
1	L5	4656	A
1	L5	4657	U
1	L5	4670	C
1	L5	4672	A
1	L5	4679	G
1	L5	4684	A
1	L5	4687	A
1	L5	4694	G
1	L5	4695	C
1	L5	4700	A
1	L5	4708	A
1	L5	4709	U
1	L5	4719	G
1	L5	4720	C
1	L5	4731	G
1	L5	4734	A
1	L5	4741	C
1	L5	4742	G
1	L5	4745	G
1	L5	4750	G
1	L5	4754	G
1	L5	4757	C
1	L5	4759	C
1	L5	4761	G
1	L5	4764	A
1	L5	4765	G
1	L5	4770	U
1	L5	4771	C
1	L5	4772	C
1	L5	4773	C
1	L5	4775	C
1	L5	4776	G

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Mol	Chain	Res	Type
1	L5	4859	C
1	L5	4860	G
1	L5	4870	G
1	L5	4871	C
1	L5	4875	G
1	L5	4882	U
1	L5	4883	C
1	L5	4888	U
1	L5	4889	G
1	L5	4891	G
1	L5	4895	C
1	L5	4896	G
1	L5	4900	C
1	L5	4901	G
1	L5	4902	C
1	L5	4910	G
1	L5	4911	A
1	L5	4912	G
1	L5	4914	C
1	L5	4922	C
1	L5	4923	C
1	L5	4925	U
1	L5	4926	C
1	L5	4934	A
1	L5	4937	C
1	L5	4940	C
1	L5	4941	G
1	L5	4943	A
1	L5	4947	U
1	L5	4951	G
1	L5	4960	G
1	L5	4961	G
1	L5	4966	A
1	L5	4976	U
1	L5	4979	A
1	L5	4988	U
1	L5	4989	U
1	L5	4991	U
1	L5	5007	A
1	L5	5013	C
1	L5	5014	A
1	L5	5017	G

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Mol	Chain	Res	Type
1	L5	5022	U
1	L5	5024	C
1	L5	5025	C
1	L5	5026	U
1	L5	5027	C
1	L5	5028	G
1	L5	5029	C
1	L5	5034	A
1	L5	5041	G
1	L5	5047	C
1	L5	5050	C
1	L5	5054	C
1	L5	5055	G
1	L5	5061	A
1	L5	5069	U
2	L7	33	U
2	L7	38	U
2	L7	53	U
2	L7	64	G
2	L7	97	G
2	L7	100	A
2	L7	106	G
2	L7	110	G
2	L7	111	C
3	L8	25	G
3	L8	34	U
3	L8	35	C
3	L8	39	G
3	L8	48	A
3	L8	59	A
3	L8	62	A
3	L8	63	U
3	L8	68	G
3	L8	82	A
3	L8	83	C
3	L8	84	A
3	L8	85	U
3	L8	86	U
3	L8	87	G
3	L8	94	G
3	L8	103	A
3	L8	105	C

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Mol	Chain	Res	Type
3	L8	110	U
3	L8	112	G
3	L8	114	G
3	L8	123	U
3	L8	124	U
3	L8	125	C
3	L8	126	C
3	L8	127	U
49	S2	2	A
49	S2	5	U
49	S2	23	G
49	S2	25	A
49	S2	33	G
49	S2	41	G
49	S2	42	A
49	S2	44	U
49	S2	46	A
49	S2	56	G
49	S2	58	C
49	S2	59	U
49	S2	62	G
49	S2	64	A
49	S2	65	C
49	S2	67	C
49	S2	68	A
49	S2	72	C
49	S2	73	C
49	S2	74	G
49	S2	76	U
49	S2	103	A
49	S2	113	G
49	S2	115	U
49	S2	116	U
49	S2	120	U
49	S2	121	U
49	S2	123	G
49	S2	126	G
49	S2	128	U
49	S2	129	C
49	S2	130	G
49	S2	139	C
49	S2	142	C

Continued on next page...

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Mol	Chain	Res	Type
49	S2	143	U
49	S2	145	G
49	S2	147	A
49	S2	149	A
49	S2	160	U
49	S2	162	C
49	S2	163	U
49	S2	166	A
49	S2	175	A
49	S2	190	G
49	S2	198	U
49	S2	199	C
49	S2	200	G
49	S2	203	G
49	S2	204	G
49	S2	207	G
49	S2	211	G
49	S2	213	G
49	S2	214	U
49	S2	219	U
49	S2	292	A
49	S2	295	C
49	S2	302	A
49	S2	305	U
49	S2	306	C
49	S2	307	G
49	S2	308	G
49	S2	309	G
49	S2	312	G
49	S2	313	A
49	S2	318	A
49	S2	319	C
49	S2	322	C
49	S2	323	C
49	S2	324	C
49	S2	325	C
49	S2	326	C
49	S2	327	G
49	S2	328	U
49	S2	329	G
49	S2	335	G
49	S2	339	A

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Mol	Chain	Res	Type
49	S2	347	G
49	S2	360	A
49	S2	362	C
49	S2	364	A
49	S2	368	U
49	S2	370	G
49	S2	385	G
49	S2	386	C
49	S2	399	C
49	S2	400	C
49	S2	408	A
49	S2	409	C
49	S2	438	G
49	S2	448	A
49	S2	449	A
49	S2	450	C
49	S2	452	G
49	S2	459	C
49	S2	464	A
49	S2	465	A
49	S2	466	G
49	S2	471	G
49	S2	472	C
49	S2	473	A
49	S2	474	G
49	S2	482	G
49	S2	487	U
49	S2	489	A
49	S2	492	C
49	S2	493	A
49	S2	496	C
49	S2	503	C
49	S2	504	G
49	S2	506	G
49	S2	507	G
49	S2	516	A
49	S2	523	A
49	S2	530	U
49	S2	540	U
49	S2	541	U
49	S2	542	U
49	S2	545	A

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Mol	Chain	Res	Type
49	S2	546	G
49	S2	547	G
49	S2	548	C
49	S2	554	A
49	S2	560	A
49	S2	563	G
49	S2	564	A
49	S2	576	A
49	S2	583	C
49	S2	585	C
49	S2	589	G
49	S2	590	A
49	S2	591	U
49	S2	592	C
49	S2	597	G
49	S2	600	G
49	S2	601	G
49	S2	604	A
49	S2	608	C
49	S2	613	G
49	S2	614	C
49	S2	617	G
49	S2	623	G
49	S2	627	U
49	S2	628	A
49	S2	629	A
49	S2	632	C
49	S2	643	A
49	S2	644	G
49	S2	660	C
49	S2	663	C
49	S2	664	A
49	S2	668	A
49	S2	669	A
49	S2	671	A
49	S2	672	A
49	S2	673	G
49	S2	678	U
49	S2	680	G
49	S2	683	G
49	S2	684	G
49	S2	687	C

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Mol	Chain	Res	Type
49	S2	688	U
49	S2	689	U
49	S2	690	G
49	S2	692	G
49	S2	693	A
49	S2	695	C
49	S2	696	G
49	S2	697	G
49	S2	698	G
49	S2	732	U
49	S2	733	C
49	S2	736	C
49	S2	738	C
49	S2	749	U
49	S2	751	G
49	S2	752	G
49	S2	753	C
49	S2	788	G
49	S2	791	C
49	S2	792	C
49	S2	798	A
49	S2	799	U
49	S2	809	A
49	S2	821	G
49	S2	822	U
49	S2	827	A
49	S2	830	A
49	S2	834	C
49	S2	835	C
49	S2	836	G
49	S2	837	A
49	S2	838	G
49	S2	839	C
49	S2	840	C
49	S2	841	G
49	S2	842	C
49	S2	845	G
49	S2	847	A
49	S2	853	C
49	S2	869	A
49	S2	870	A
49	S2	871	U

Continued on next page...

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Mol	Chain	Res	Type
49	S2	872	A
49	S2	873	G
49	S2	874	G
49	S2	877	C
49	S2	880	G
49	S2	883	U
49	S2	887	U
49	S2	888	U
49	S2	889	U
49	S2	891	G
49	S2	893	U
49	S2	896	U
49	S2	897	U
49	S2	898	U
49	S2	899	U
49	S2	900	C
49	S2	901	G
49	S2	903	A
49	S2	904	A
49	S2	906	U
49	S2	913	A
49	S2	920	A
49	S2	922	A
49	S2	933	G
49	S2	934	G
49	S2	937	C
49	S2	943	U
49	S2	955	A
49	S2	956	G
49	S2	961	G
49	S2	963	A
49	S2	968	U
49	S2	970	G
49	S2	985	G
49	S2	990	A
49	S2	992	A
49	S2	997	A
49	S2	999	G
49	S2	1001	A
49	S2	1002	U
49	S2	1011	A
49	S2	1016	U

Continued on next page...

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Mol	Chain	Res	Type
49	S2	1017	U
49	S2	1023	A
49	S2	1033	G
49	S2	1039	C
49	S2	1061	U
49	S2	1062	A
49	S2	1066	U
49	S2	1067	C
49	S2	1069	U
49	S2	1083	A
49	S2	1085	C
49	S2	1089	G
49	S2	1100	A
49	S2	1109	C
49	S2	1114	U
49	S2	1115	U
49	S2	1116	C
49	S2	1118	C
49	S2	1119	A
49	S2	1121	G
49	S2	1133	A
49	S2	1138	C
49	S2	1140	G
49	S2	1150	A
49	S2	1153	C
49	S2	1154	U
49	S2	1195	A
49	S2	1203	G
49	S2	1209	A
49	S2	1212	G
49	S2	1215	C
49	S2	1216	C
49	S2	1217	A
49	S2	1221	G
49	S2	1224	G
49	S2	1227	G
49	S2	1231	C
49	S2	1242	U
49	S2	1243	U
49	S2	1251	A
49	S2	1253	A
49	S2	1256	G

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Mol	Chain	Res	Type
49	S2	1257	G
49	S2	1259	A
49	S2	1271	C
49	S2	1274	G
49	S2	1275	G
49	S2	1283	C
49	S2	1284	A
49	S2	1286	G
49	S2	1294	G
49	S2	1295	A
49	S2	1301	A
49	S2	1302	G
49	S2	1303	C
49	S2	1306	U
49	S2	1308	U
49	S2	1312	G
49	S2	1314	U
49	S2	1330	G
49	S2	1341	C
49	S2	1342	U
49	S2	1348	G
49	S2	1353	A
49	S2	1358	U
49	S2	1363	C
49	S2	1371	U
49	S2	1372	U
49	S2	1373	C
49	S2	1378	A
49	S2	1382	A
49	S2	1396	A
49	S2	1397	U
49	S2	1402	A
49	S2	1404	U
49	S2	1412	C
49	S2	1415	C
49	S2	1416	C
49	S2	1421	A
49	S2	1423	C
49	S2	1424	G
49	S2	1431	G
49	S2	1433	C
49	S2	1434	C

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Mol	Chain	Res	Type
49	S2	1435	C
49	S2	1436	C
49	S2	1438	A
49	S2	1449	G
49	S2	1452	A
49	S2	1454	A
49	S2	1455	A
49	S2	1463	U
49	S2	1464	C
49	S2	1478	U
49	S2	1480	A
49	S2	1486	A
49	S2	1489	A
49	S2	1490	G
49	S2	1497	G
49	S2	1498	A
49	S2	1506	A
49	S2	1507	G
49	S2	1508	A
49	S2	1520	G
49	S2	1521	C
49	S2	1523	C
49	S2	1531	A
49	S2	1533	A
49	S2	1534	C
49	S2	1539	U
49	S2	1544	C
49	S2	1549	U
49	S2	1551	U
49	S2	1552	G
49	S2	1553	C
49	S2	1556	A
49	S2	1558	C
49	S2	1560	U
49	S2	1570	G
49	S2	1573	G
49	S2	1575	G
49	S2	1579	A
49	S2	1580	A
49	S2	1585	U
49	S2	1587	G
49	S2	1588	A

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Mol	Chain	Res	Type
49	S2	1594	A
49	S2	1595	U
49	S2	1598	G
49	S2	1600	G
49	S2	1601	A
49	S2	1603	G
49	S2	1606	G
49	S2	1612	G
49	S2	1621	U
49	S2	1623	A
49	S2	1629	C
49	S2	1632	G
49	S2	1634	A
49	S2	1643	U
49	S2	1644	C
49	S2	1646	C
49	S2	1647	A
49	S2	1648	G
49	S2	1649	U
49	S2	1653	U
49	S2	1654	G
49	S2	1663	A
49	S2	1664	A
49	S2	1665	G
49	S2	1671	G
49	S2	1677	U
49	S2	1680	G
49	S2	1691	U
49	S2	1696	C
49	S2	1698	C
49	S2	1699	A
49	S2	1700	C
49	S2	1701	C
49	S2	1715	A
49	S2	1719	A
49	S2	1721	U
49	S2	1722	G
49	S2	1742	C
49	S2	1744	G
49	S2	1745	A
49	S2	1748	G
49	S2	1750	C

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Mol	Chain	Res	Type
49	S2	1752	C
49	S2	1753	C
49	S2	1754	G
49	S2	1757	G
49	S2	1758	G
49	S2	1759	G
49	S2	1760	G
49	S2	1761	U
49	S2	1771	G
49	S2	1772	C
49	S2	1773	C
49	S2	1774	C
49	S2	1775	U
49	S2	1776	G
49	S2	1777	G
49	S2	1781	A
49	S2	1783	C
49	S2	1784	G
49	S2	1786	U
49	S2	1787	G
49	S2	1788	A
49	S2	1798	C
49	S2	1819	A
49	S2	1822	A
49	S2	1823	A
49	S2	1824	A
49	S2	1825	A
49	S2	1826	G
49	S2	1829	G
49	S2	1831	A
49	S2	1835	A
49	S2	1838	U
49	S2	1849	G
49	S2	1851	A
49	S2	1852	C
49	S2	1861	G
49	S2	1862	G
49	S2	1863	A
49	S2	1864	U
49	S2	1865	C
85	CC	7	G
85	CC	8	U

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Mol	Chain	Res	Type
85	CC	9	G
85	CC	17	G
85	CC	18	G
85	CC	20	A
85	CC	29	G
85	CC	37	A
85	CC	43	A
85	CC	46	U
85	CC	47	C
85	CC	57	A
85	CC	58	A
85	CC	60	C
85	CC	63	U
85	CC	73	C
85	CC	75	A

All (26) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	L5	184	U
1	L5	186	G
1	L5	406	C
1	L5	493	G
1	L5	504	G
1	L5	914	U
1	L5	1633	G
1	L5	1977	C
1	L5	2033	A
1	L5	2416	G
1	L5	2475	G
1	L5	2675	G
1	L5	2760	G
1	L5	2786	C
1	L5	3614	G
1	L5	3673	C
1	L5	3735	G
1	L5	4378	A
1	L5	4699	U
1	L5	4913	G
49	S2	112	U
49	S2	291	G
49	S2	563	G

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Mol	Chain	Res	Type
49	S2	688	U
49	S2	1434	C
85	CC	73	C

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
41	MLZ	Lm	98	41	8,9,10	0.78	0	4,9,11	0.61	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
41	MLZ	Lm	98	41	-	1/7/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
41	Lm	98	MLZ	CD-CE-NZ-CM

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 271 ligands modelled in this entry, 263 are monoatomic - leaving 8 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	T1C	L5	5313	87	44,45,45	1.18	4 (9%)	53,72,72	1.11	3 (5%)
88	T1C	L5	5312	-	44,45,45	1.19	4 (9%)	53,72,72	1.06	3 (5%)
88	T1C	L5	5315	-	44,45,45	1.21	4 (9%)	53,72,72	1.40	7 (13%)
88	T1C	CC	101	-	44,45,45	1.18	4 (9%)	53,72,72	1.05	3 (5%)
88	T1C	L5	5314	87	44,45,45	1.19	4 (9%)	53,72,72	1.11	5 (9%)
88	T1C	L5	5316	-	44,45,45	1.24	4 (9%)	53,72,72	1.61	12 (22%)
88	T1C	L5	5317	-	44,45,45	1.20	4 (9%)	53,72,72	1.15	6 (11%)
88	T1C	L5	5311	-	44,45,45	1.18	4 (9%)	53,72,72	1.06	2 (3%)

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	T1C	L5	5313	87	-	9/22/80/80	0/4/4/4
88	T1C	L5	5312	-	-	16/22/80/80	0/4/4/4
88	T1C	L5	5315	-	-	16/22/80/80	0/4/4/4
88	T1C	CC	101	-	-	6/22/80/80	0/4/4/4
88	T1C	L5	5314	87	-	13/22/80/80	0/4/4/4
88	T1C	L5	5316	-	-	9/22/80/80	0/4/4/4
88	T1C	L5	5317	-	-	15/22/80/80	0/4/4/4
88	T1C	L5	5311	-	-	12/22/80/80	0/4/4/4

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
88	L5	5317	T1C	C21-N21	5.52	1.48	1.33
88	L5	5315	T1C	C21-N21	5.49	1.48	1.33
88	L5	5313	T1C	C21-N21	5.46	1.47	1.33
88	L5	5311	T1C	C21-N21	5.45	1.47	1.33
88	L5	5312	T1C	C21-N21	5.44	1.47	1.33
88	L5	5316	T1C	C21-N21	5.43	1.47	1.33
88	CC	101	T1C	C21-N21	5.42	1.47	1.33
88	L5	5314	T1C	C21-N21	5.36	1.47	1.33
88	L5	5312	T1C	C4-N4	2.57	1.53	1.47
88	L5	5314	T1C	C4-N4	2.54	1.53	1.47
88	L5	5317	T1C	C4-N4	2.52	1.53	1.47
88	L5	5316	T1C	C7-N7	2.43	1.49	1.42
88	CC	101	T1C	C4-N4	2.33	1.52	1.47
88	L5	5311	T1C	C4-N4	2.31	1.52	1.47
88	L5	5313	T1C	O11-C11	2.28	1.28	1.23
88	L5	5313	T1C	C4-N4	2.24	1.52	1.47
88	L5	5316	T1C	C4-N4	2.23	1.52	1.47
88	L5	5314	T1C	O11-C11	2.22	1.27	1.23
88	L5	5317	T1C	O11-C11	2.21	1.27	1.23
88	CC	101	T1C	O11-C11	2.20	1.27	1.23
88	L5	5315	T1C	C4-N4	2.18	1.52	1.47
88	L5	5316	T1C	O11-C11	2.17	1.27	1.23
88	L5	5312	T1C	O11-C11	2.17	1.27	1.23
88	L5	5315	T1C	O11-C11	2.17	1.27	1.23
88	L5	5311	T1C	O11-C11	2.14	1.27	1.23
88	L5	5317	T1C	C7-N7	2.13	1.48	1.42
88	CC	101	T1C	C7-N7	2.13	1.48	1.42
88	L5	5315	T1C	C7-N7	2.12	1.48	1.42
88	L5	5312	T1C	C7-N7	2.11	1.48	1.42
88	L5	5311	T1C	C7-N7	2.11	1.48	1.42
88	L5	5313	T1C	C7-N7	2.06	1.48	1.42
88	L5	5314	T1C	C7-N7	2.05	1.48	1.42

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
88	L5	5316	T1C	C1C-C1-C2	4.80	123.37	115.75
88	L5	5315	T1C	C11-C1B-C12	4.69	122.51	118.80
88	L5	5313	T1C	C11-C1B-C12	4.66	122.49	118.80
88	L5	5316	T1C	C51-C5-C41	4.60	118.57	110.49
88	L5	5316	T1C	C1C-C41-C4	4.54	117.85	111.64
88	L5	5315	T1C	C1C-C41-C4	4.53	117.83	111.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
88	L5	5312	T1C	C1C-C1-C2	3.92	121.98	115.75
88	L5	5311	T1C	C1C-C1-C2	3.88	121.91	115.75
88	L5	5311	T1C	C11-C1B-C12	3.73	121.75	118.80
88	CC	101	T1C	C11-C1B-C12	3.67	121.70	118.80
88	L5	5315	T1C	C1C-C1-C2	3.31	121.00	115.75
88	L5	5313	T1C	C1-C1C-C12	3.23	113.67	109.88
88	L5	5317	T1C	C1C-C1-C2	3.13	120.73	115.75
88	CC	101	T1C	C1-C1C-C12	3.02	113.43	109.88
88	L5	5316	T1C	O12-C12-C1B	-2.99	119.81	123.90
88	L5	5314	T1C	C1C-C1-C2	2.95	120.44	115.75
88	L5	5317	T1C	C11-C1B-C12	2.83	121.04	118.80
88	L5	5315	T1C	C51-C5-C41	2.82	115.45	110.49
88	L5	5316	T1C	O1C-C1C-C12	-2.70	105.82	110.14
88	L5	5316	T1C	C61-C7-N7	2.70	122.21	118.91
88	L5	5316	T1C	C11-C1B-C12	2.65	120.89	118.80
88	L5	5314	T1C	O1C-C1C-C12	-2.63	105.94	110.14
88	L5	5312	T1C	C51-C5-C41	2.54	114.95	110.49
88	L5	5315	T1C	O1C-C1C-C12	-2.48	106.17	110.14
88	L5	5312	T1C	C11-C1B-C12	2.42	120.72	118.80
88	L5	5316	T1C	C1C-C12-C1B	2.42	125.52	123.06
88	L5	5316	T1C	C8-C7-N7	-2.39	117.71	120.91
88	CC	101	T1C	O1C-C1C-C12	-2.38	106.33	110.14
88	L5	5314	T1C	C1C-C41-C4	2.38	114.89	111.64
88	L5	5314	T1C	C11-C1B-C12	2.32	120.64	118.80
88	L5	5317	T1C	O12-C12-C1B	-2.26	120.80	123.90
88	L5	5317	T1C	C1C-C41-C4	2.25	114.72	111.64
88	L5	5315	T1C	C1C-C12-C1B	-2.22	120.81	123.06
88	L5	5317	T1C	C1-C1C-C12	2.22	112.48	109.88
88	L5	5316	T1C	C72-N7-C71	-2.21	109.02	116.12
88	L5	5316	T1C	O1C-C1C-C41	-2.20	107.39	110.09
88	L5	5317	T1C	O1C-C1C-C12	-2.18	106.66	110.14
88	L5	5315	T1C	O1C-C1C-C41	-2.17	107.43	110.09
88	L5	5313	T1C	O1C-C1C-C12	-2.10	106.78	110.14
88	L5	5316	T1C	O1-C1-C2	-2.07	119.21	123.55
88	L5	5314	T1C	C6-C61-C7	-2.04	118.59	123.52

There are no chirality outliers.

All (96) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
88	L5	5311	T1C	C3-C2-C21-O21
88	L5	5311	T1C	C3-C2-C21-N21

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Mol	Chain	Res	Type	Atoms
88	L5	5311	T1C	C1-C2-C21-O21
88	L5	5311	T1C	C1-C2-C21-N21
88	L5	5312	T1C	C91-C92-N92-C93
88	L5	5312	T1C	C92-C91-N9-C9
88	L5	5312	T1C	C3-C4-N4-C43
88	L5	5312	T1C	C3-C4-N4-C42
88	L5	5312	T1C	C3-C2-C21-O21
88	L5	5312	T1C	C3-C2-C21-N21
88	L5	5312	T1C	C1-C2-C21-O21
88	L5	5313	T1C	C91-C92-N92-C93
88	L5	5313	T1C	C92-C91-N9-C9
88	L5	5313	T1C	O91-C91-N9-C9
88	L5	5313	T1C	C1-C2-C21-O21
88	L5	5313	T1C	C1-C2-C21-N21
88	L5	5314	T1C	C91-C92-N92-C93
88	L5	5314	T1C	C92-C91-N9-C9
88	L5	5314	T1C	C3-C4-N4-C43
88	L5	5314	T1C	C1-C2-C21-O21
88	L5	5314	T1C	C1-C2-C21-N21
88	L5	5315	T1C	C94-C93-N92-C92
88	L5	5315	T1C	C95-C93-N92-C92
88	L5	5315	T1C	C96-C93-N92-C92
88	L5	5315	T1C	C92-C91-N9-C9
88	L5	5315	T1C	C41-C4-N4-C43
88	L5	5315	T1C	C3-C4-N4-C43
88	L5	5315	T1C	C3-C4-N4-C42
88	L5	5315	T1C	C1-C2-C21-O21
88	L5	5315	T1C	C1-C2-C21-N21
88	L5	5316	T1C	C1-C2-C21-O21
88	L5	5316	T1C	C1-C2-C21-N21
88	L5	5317	T1C	C94-C93-N92-C92
88	L5	5317	T1C	C95-C93-N92-C92
88	L5	5317	T1C	C96-C93-N92-C92
88	L5	5317	T1C	C92-C91-N9-C9
88	L5	5317	T1C	O91-C91-N9-C9
88	L5	5317	T1C	C1-C2-C21-O21
88	L5	5317	T1C	C1-C2-C21-N21
88	CC	101	T1C	C3-C2-C21-O21
88	CC	101	T1C	C3-C2-C21-N21
88	CC	101	T1C	C1-C2-C21-O21
88	CC	101	T1C	C1-C2-C21-N21
88	L5	5315	T1C	O91-C91-N9-C9

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Mol	Chain	Res	Type	Atoms
88	L5	5314	T1C	O91-C91-N9-C9
88	L5	5312	T1C	O91-C91-N9-C9
88	L5	5311	T1C	C92-C91-N9-C9
88	L5	5317	T1C	N9-C91-C92-N92
88	L5	5317	T1C	O91-C91-C92-N92
88	L5	5315	T1C	C10-C9-N9-C91
88	L5	5311	T1C	O91-C91-N9-C9
88	L5	5311	T1C	C10-C9-N9-C91
88	L5	5315	T1C	C8-C9-N9-C91
88	CC	101	T1C	C10-C9-N9-C91
88	L5	5311	T1C	C8-C9-N9-C91
88	L5	5311	T1C	O91-C91-C92-N92
88	L5	5314	T1C	O91-C91-C92-N92
88	L5	5312	T1C	C10-C9-N9-C91
88	L5	5316	T1C	C10-C9-N9-C91
88	L5	5317	T1C	C91-C92-N92-C93
88	L5	5314	T1C	N9-C91-C92-N92
88	L5	5312	T1C	C41-C4-N4-C42
88	L5	5314	T1C	C41-C4-N4-C43
88	L5	5314	T1C	C41-C4-N4-C42
88	L5	5314	T1C	C3-C4-N4-C42
88	L5	5315	T1C	C41-C4-N4-C42
88	L5	5316	T1C	C41-C4-N4-C42
88	L5	5317	T1C	C41-C4-N4-C43
88	L5	5317	T1C	C41-C4-N4-C42
88	L5	5315	T1C	C3-C2-C21-N21
88	L5	5317	T1C	C3-C2-C21-N21
88	L5	5313	T1C	C3-C2-C21-O21
88	L5	5315	T1C	C3-C2-C21-O21
88	L5	5317	T1C	C3-C2-C21-O21
88	L5	5311	T1C	N9-C91-C92-N92
88	L5	5312	T1C	C1-C2-C21-N21
88	L5	5312	T1C	C95-C93-N92-C92
88	L5	5316	T1C	O91-C91-C92-N92
88	L5	5316	T1C	N9-C91-C92-N92
88	L5	5312	T1C	C61-C7-N7-C72
88	L5	5315	T1C	C91-C92-N92-C93
88	L5	5312	T1C	C96-C93-N92-C92
88	L5	5311	T1C	C61-C7-N7-C71
88	L5	5312	T1C	C61-C7-N7-C71
88	L5	5313	T1C	C61-C7-N7-C71
88	L5	5312	T1C	C41-C4-N4-C43

Continued on next page...

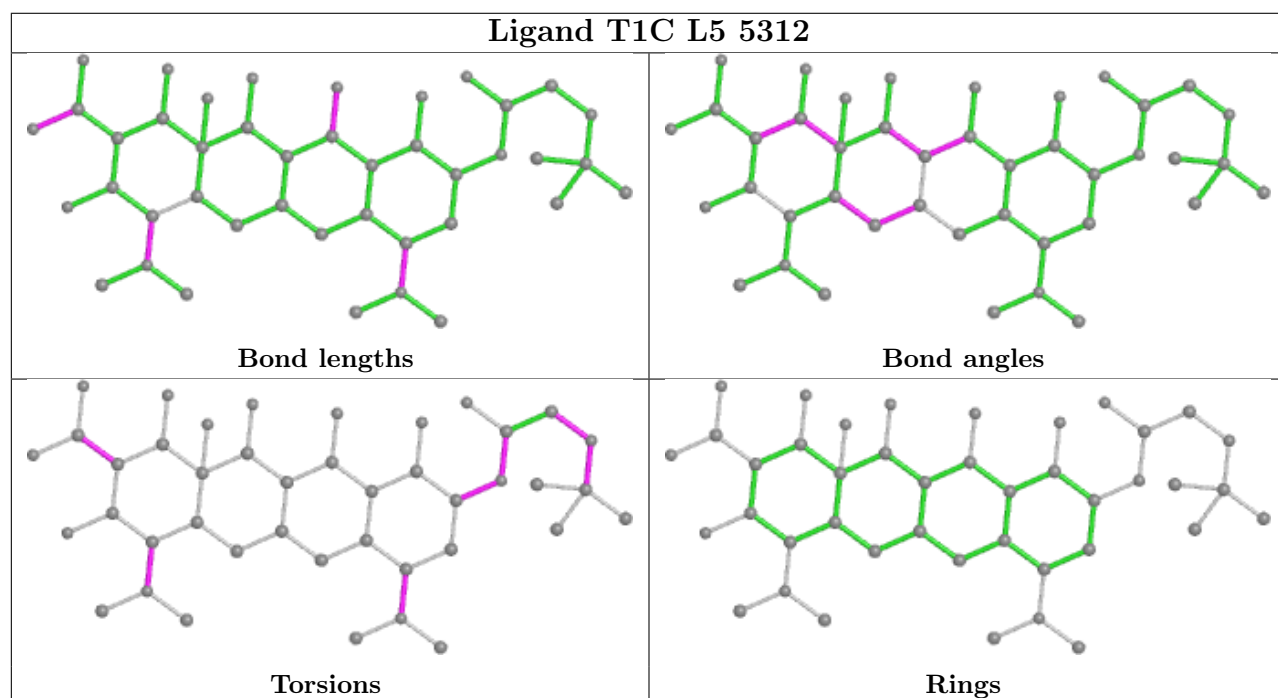
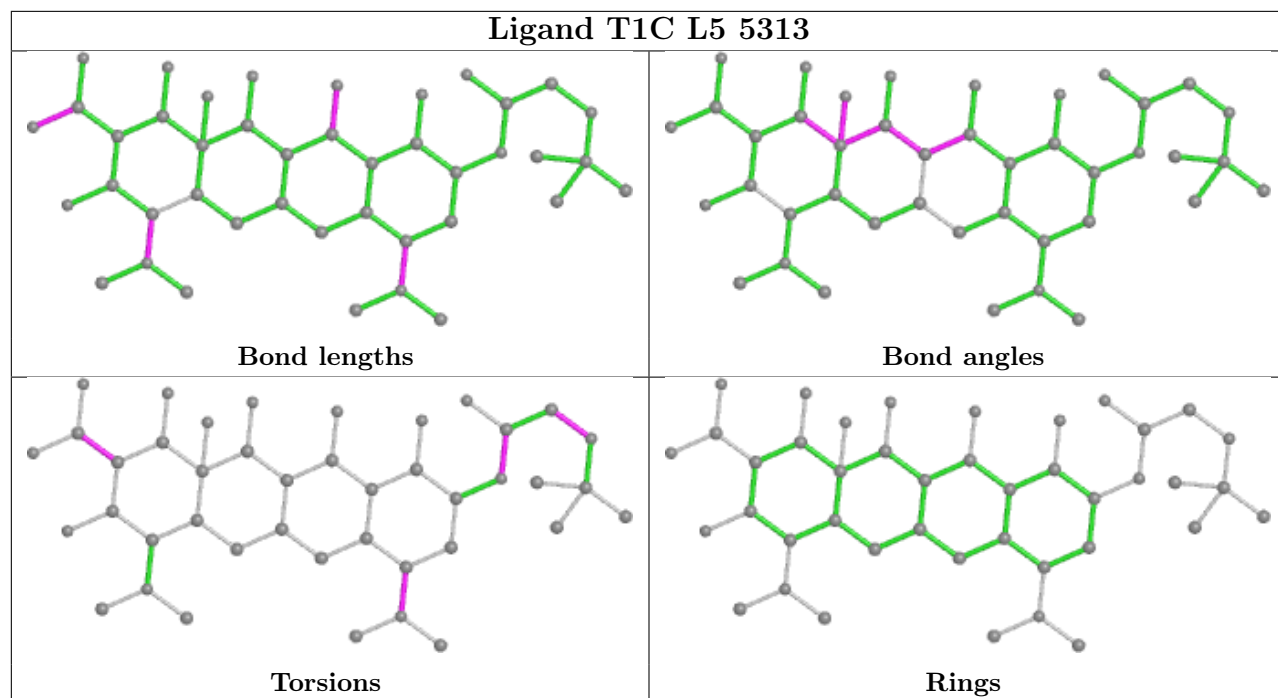
Continued from previous page...

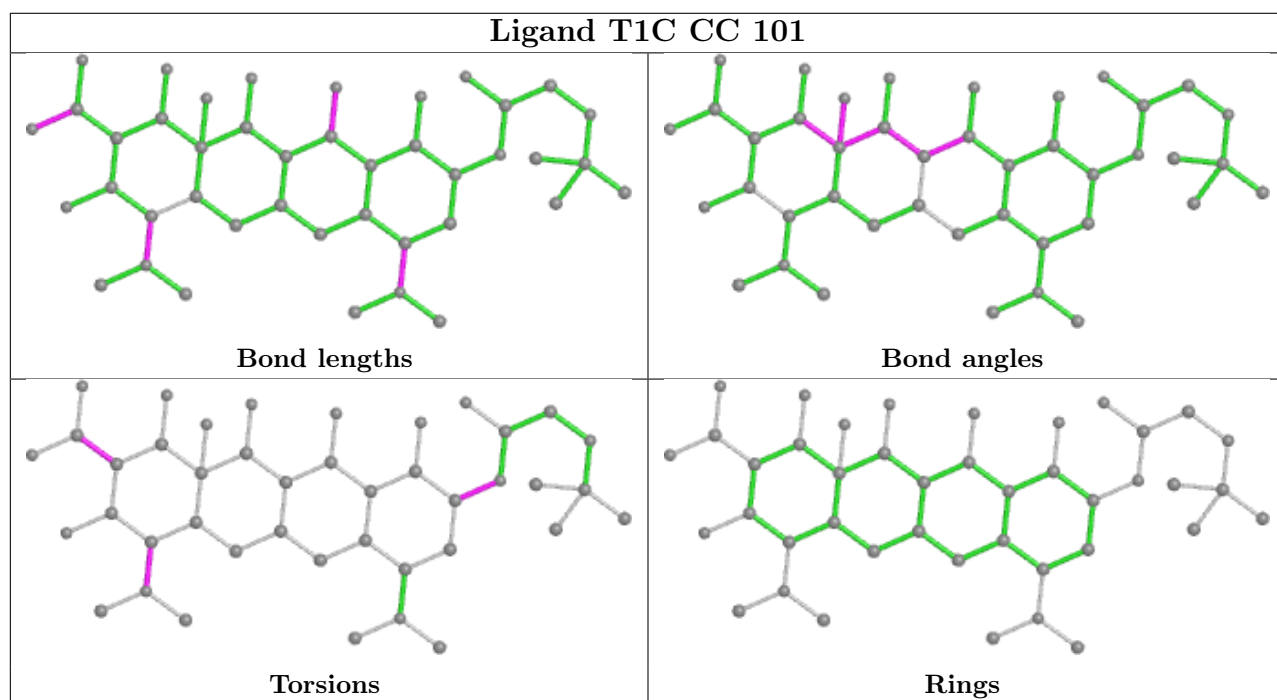
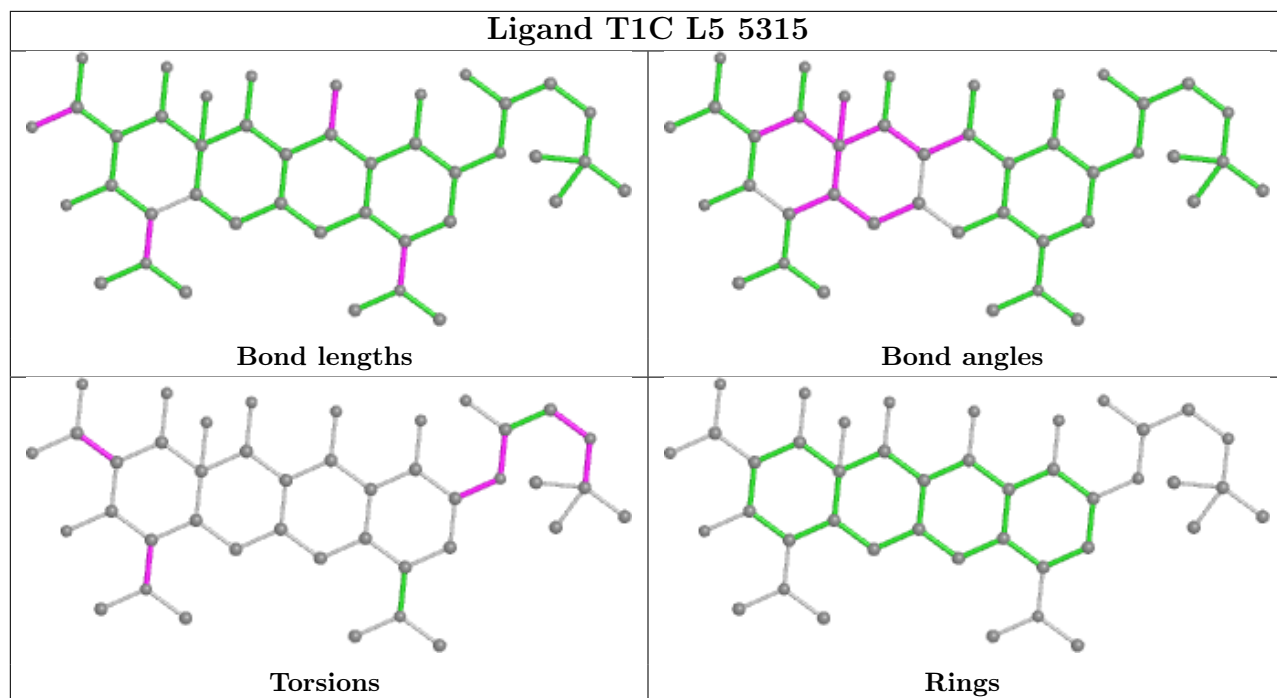
Mol	Chain	Res	Type	Atoms
88	CC	101	T1C	C41-C4-N4-C42
88	L5	5313	T1C	C61-C7-N7-C72
88	L5	5316	T1C	C8-C9-N9-C91
88	L5	5313	T1C	C3-C2-C21-N21
88	L5	5314	T1C	C3-C2-C21-N21
88	L5	5316	T1C	C3-C2-C21-N21
88	L5	5314	T1C	C3-C2-C21-O21
88	L5	5316	T1C	C3-C2-C21-O21
88	L5	5311	T1C	C61-C7-N7-C72
88	L5	5317	T1C	C61-C7-N7-C71

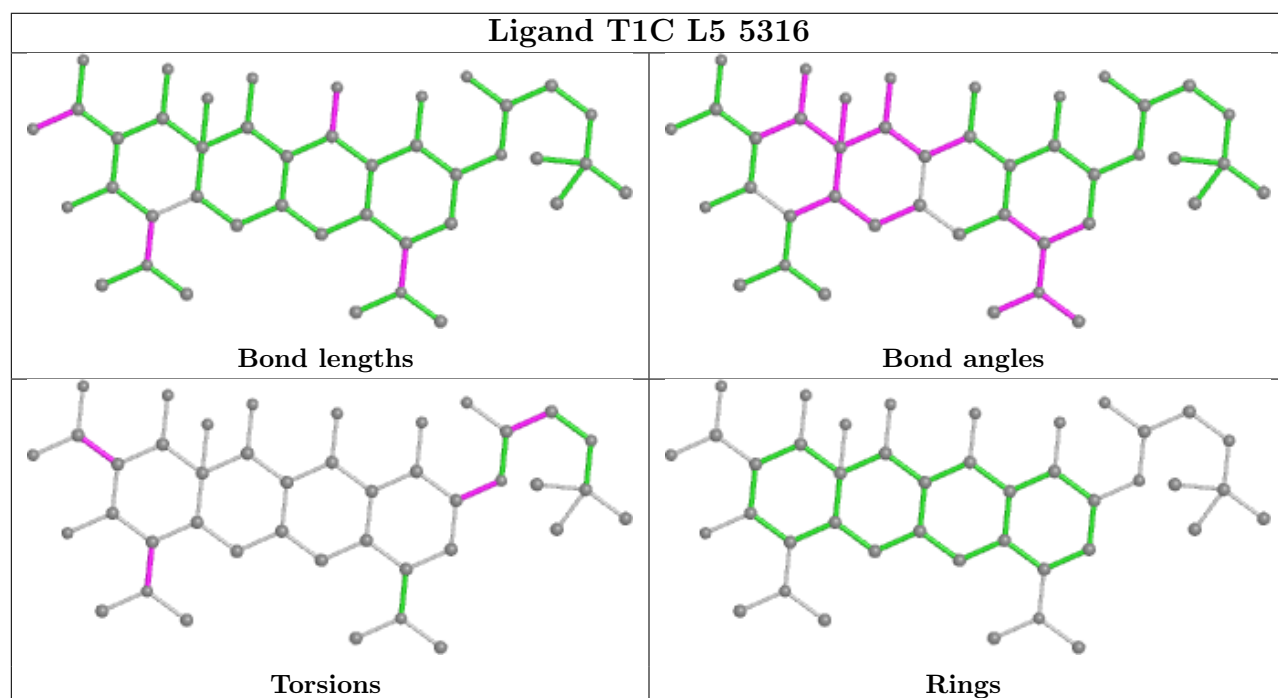
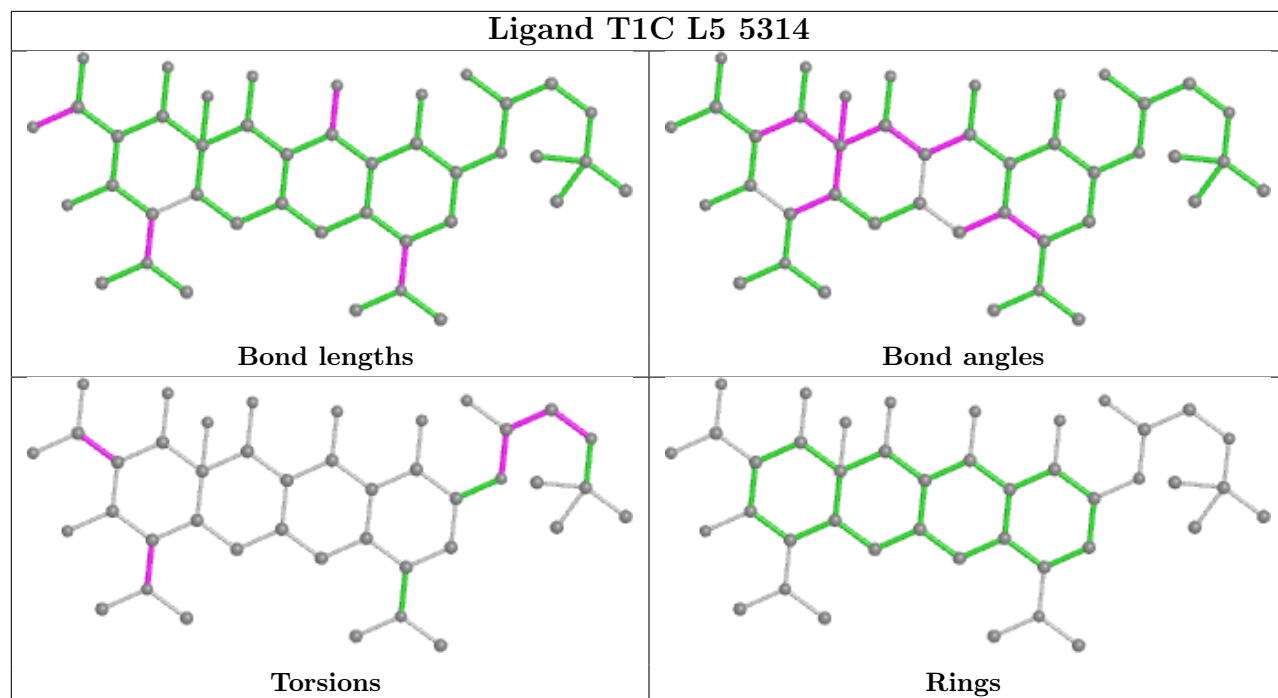
There are no ring outliers.

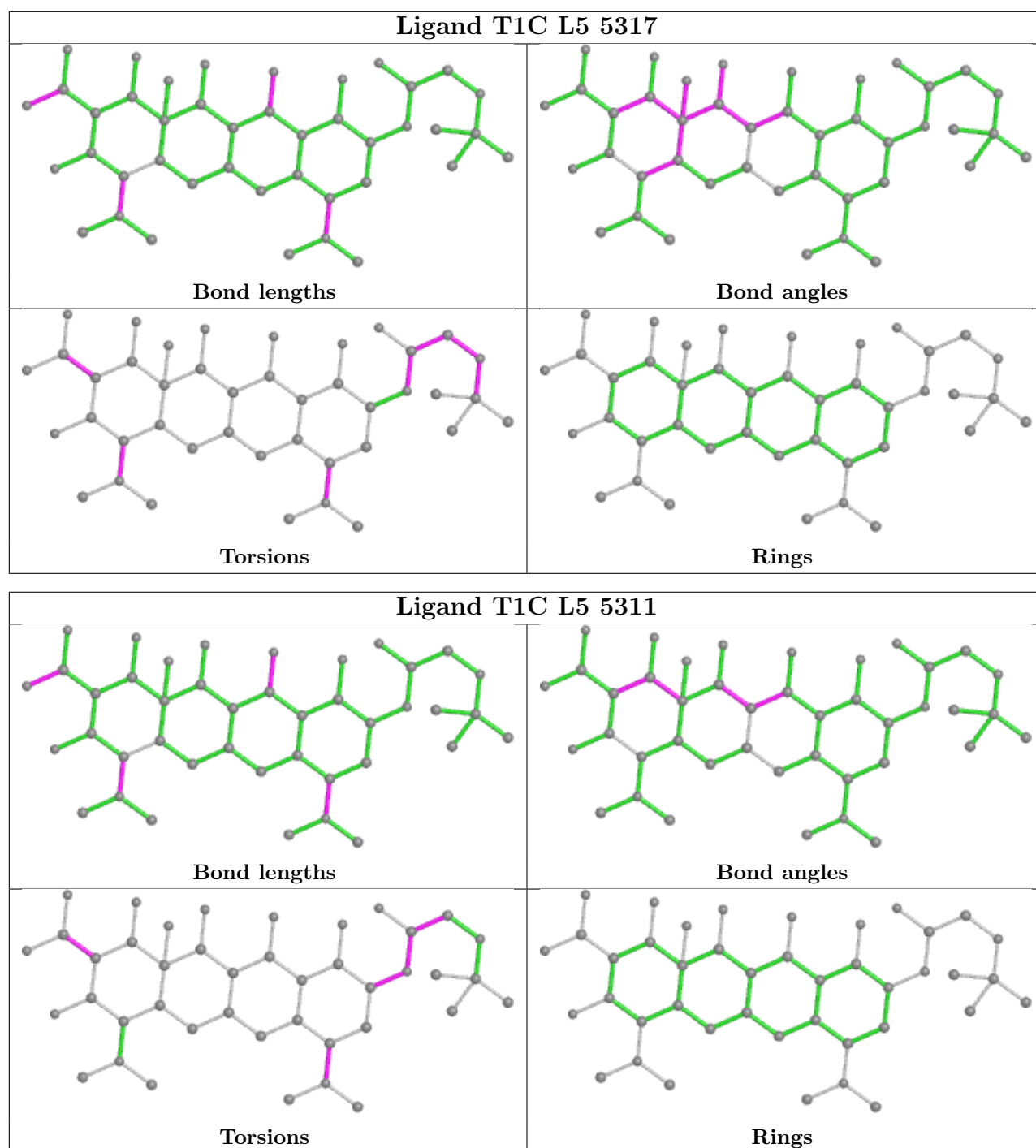
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

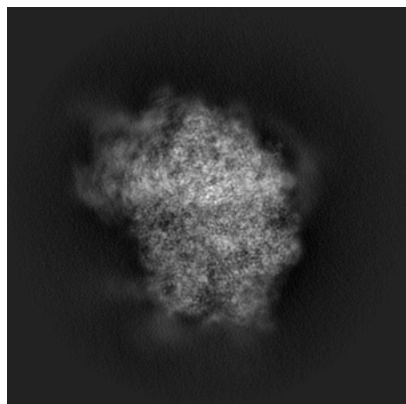
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-38630. These allow visual inspection of the internal detail of the map and identification of artifacts.

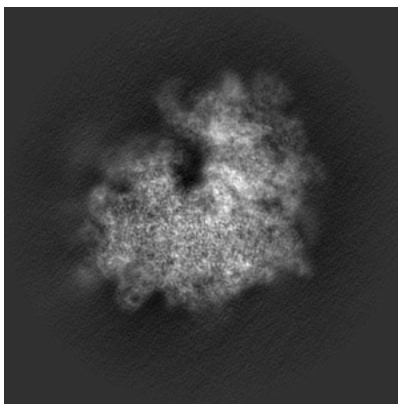
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

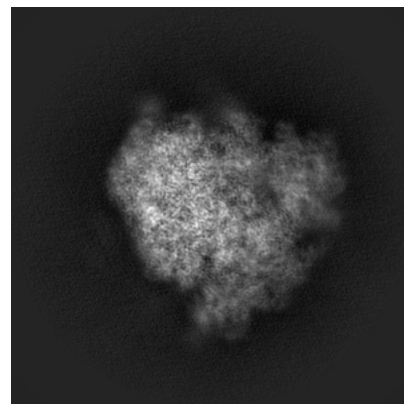
6.1.1 Primary map



X

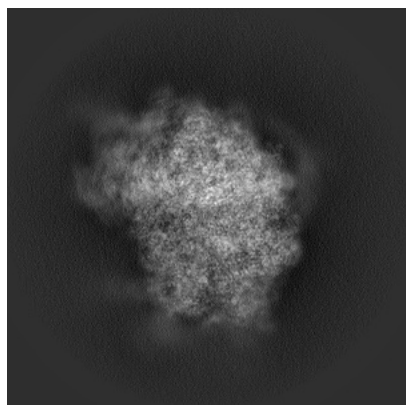


Y

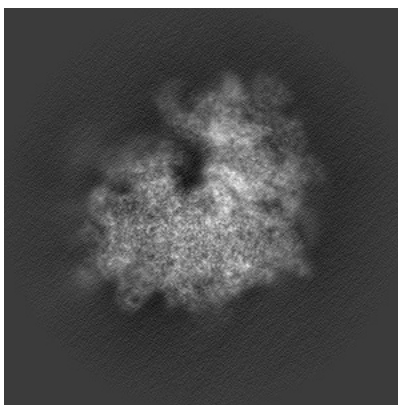


Z

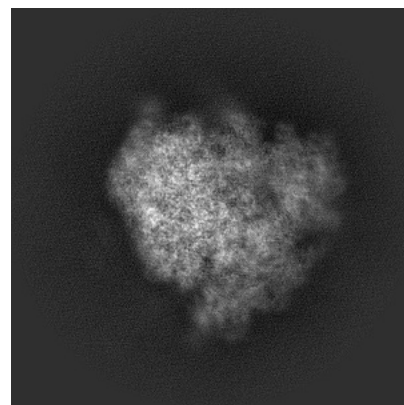
6.1.2 Raw map



X



Y

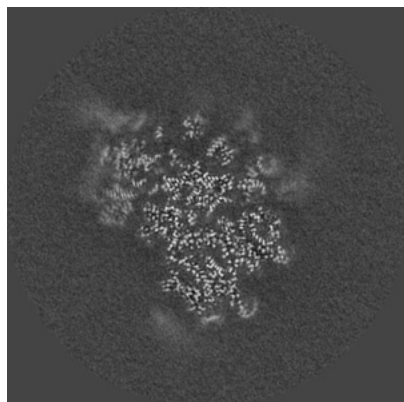


Z

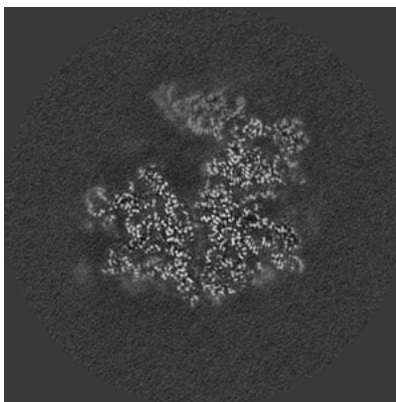
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

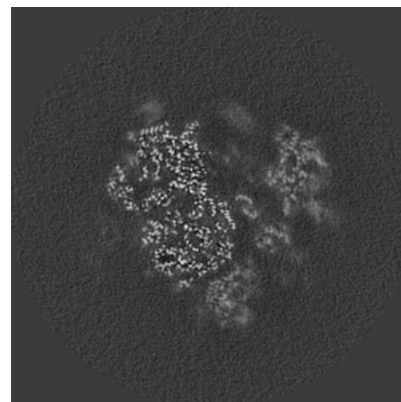
6.2.1 Primary map



X Index: 240

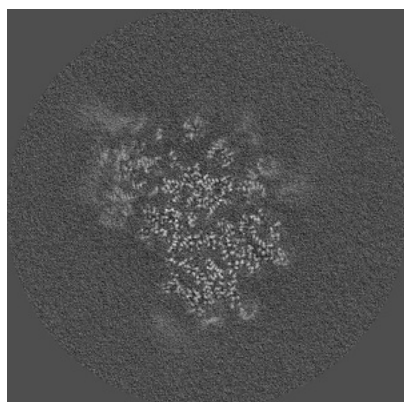


Y Index: 240

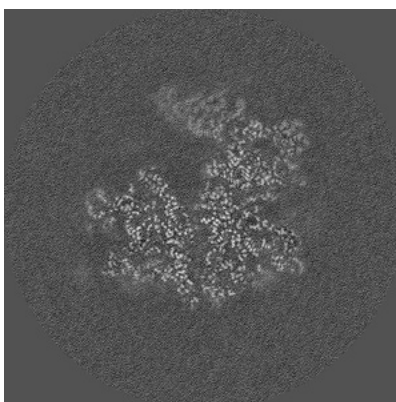


Z Index: 240

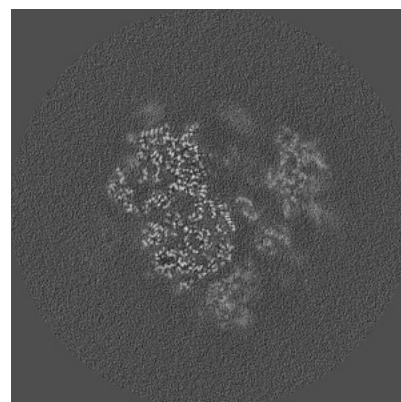
6.2.2 Raw map



X Index: 240



Y Index: 240

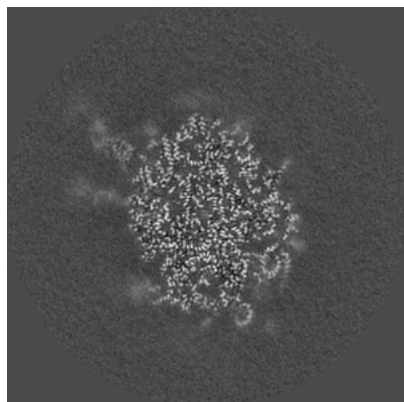


Z Index: 240

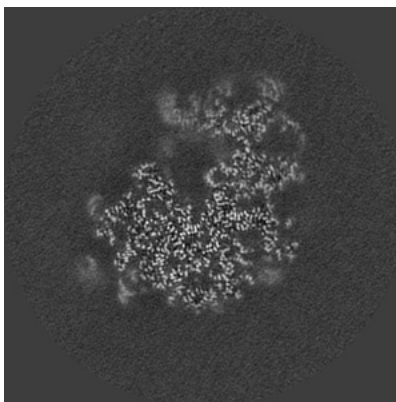
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

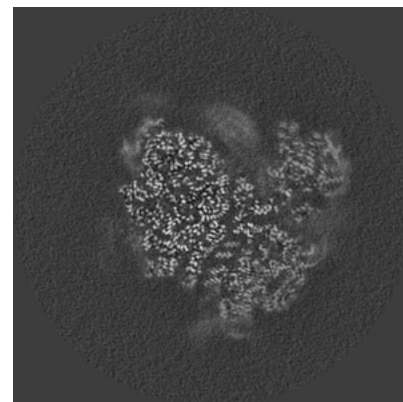
6.3.1 Primary map



X Index: 219

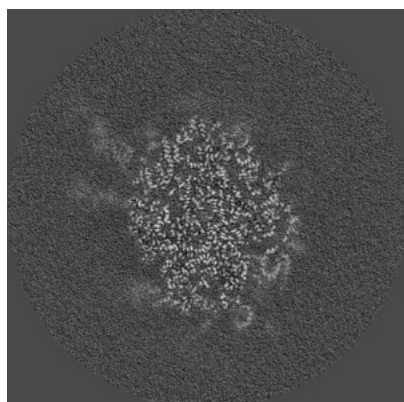


Y Index: 254

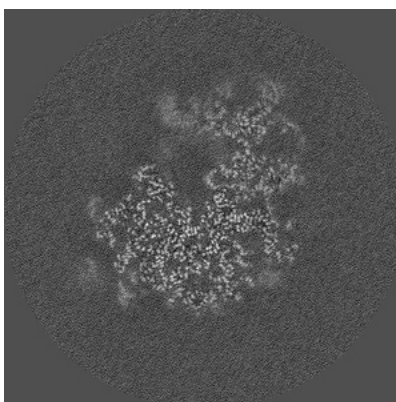


Z Index: 270

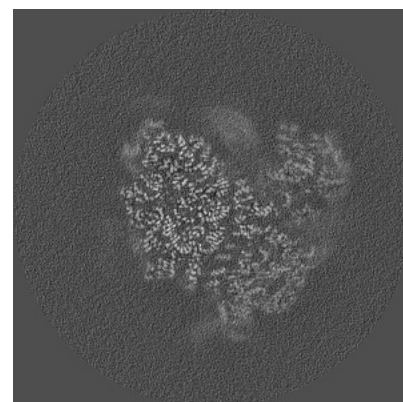
6.3.2 Raw map



X Index: 219



Y Index: 254

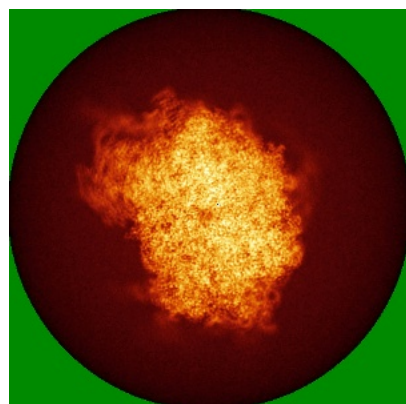


Z Index: 271

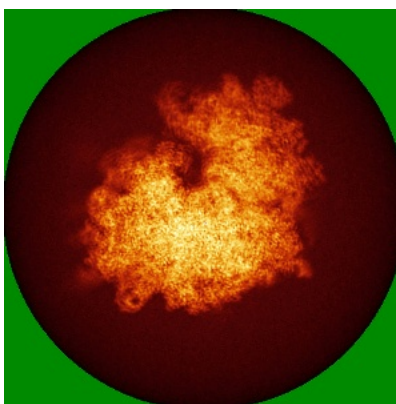
The images above show the largest variance slices of the map in three orthogonal directions.

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

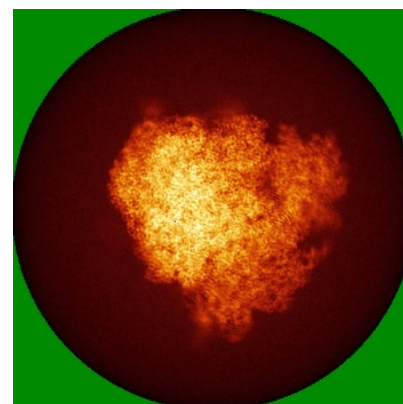
6.4.1 Primary map



X

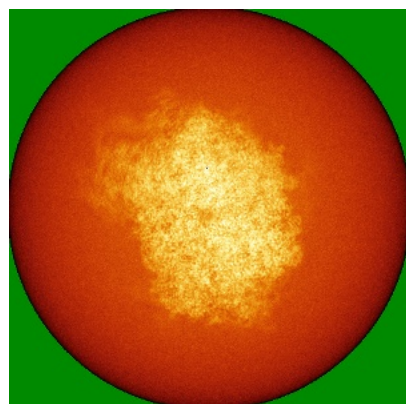


Y

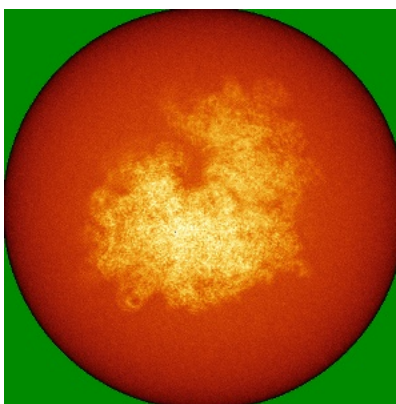


Z

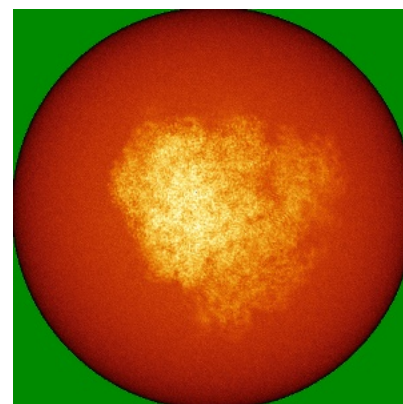
6.4.2 Raw map



X



Y

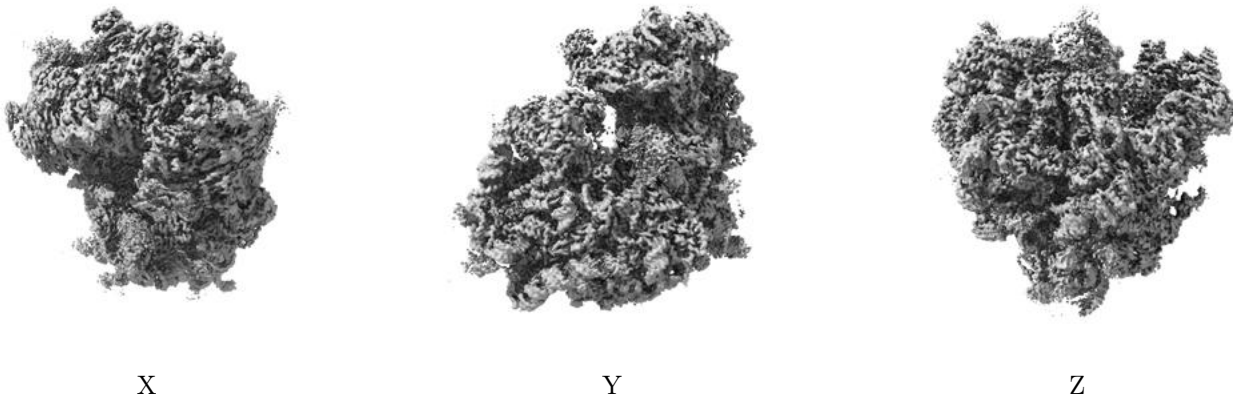


Z

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.

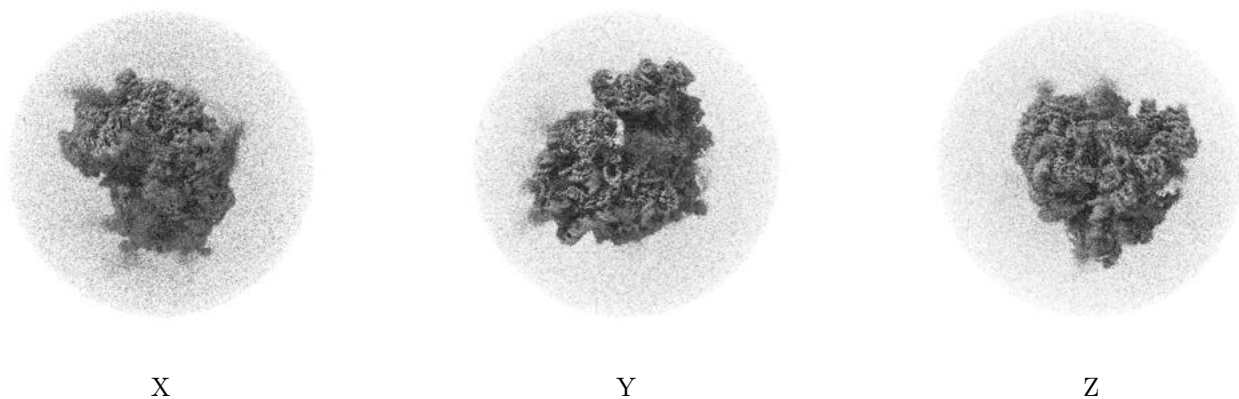
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.01. 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.

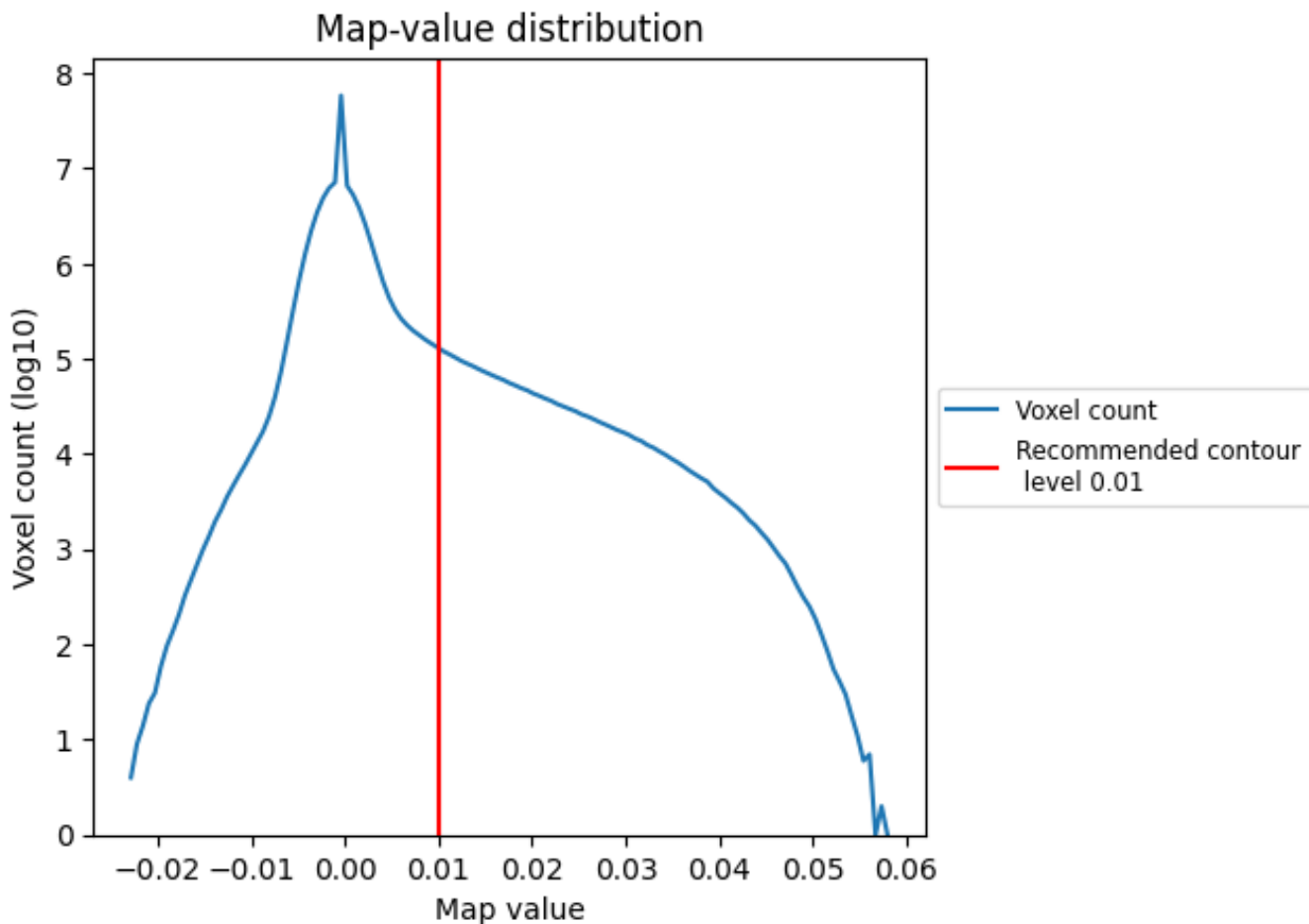
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

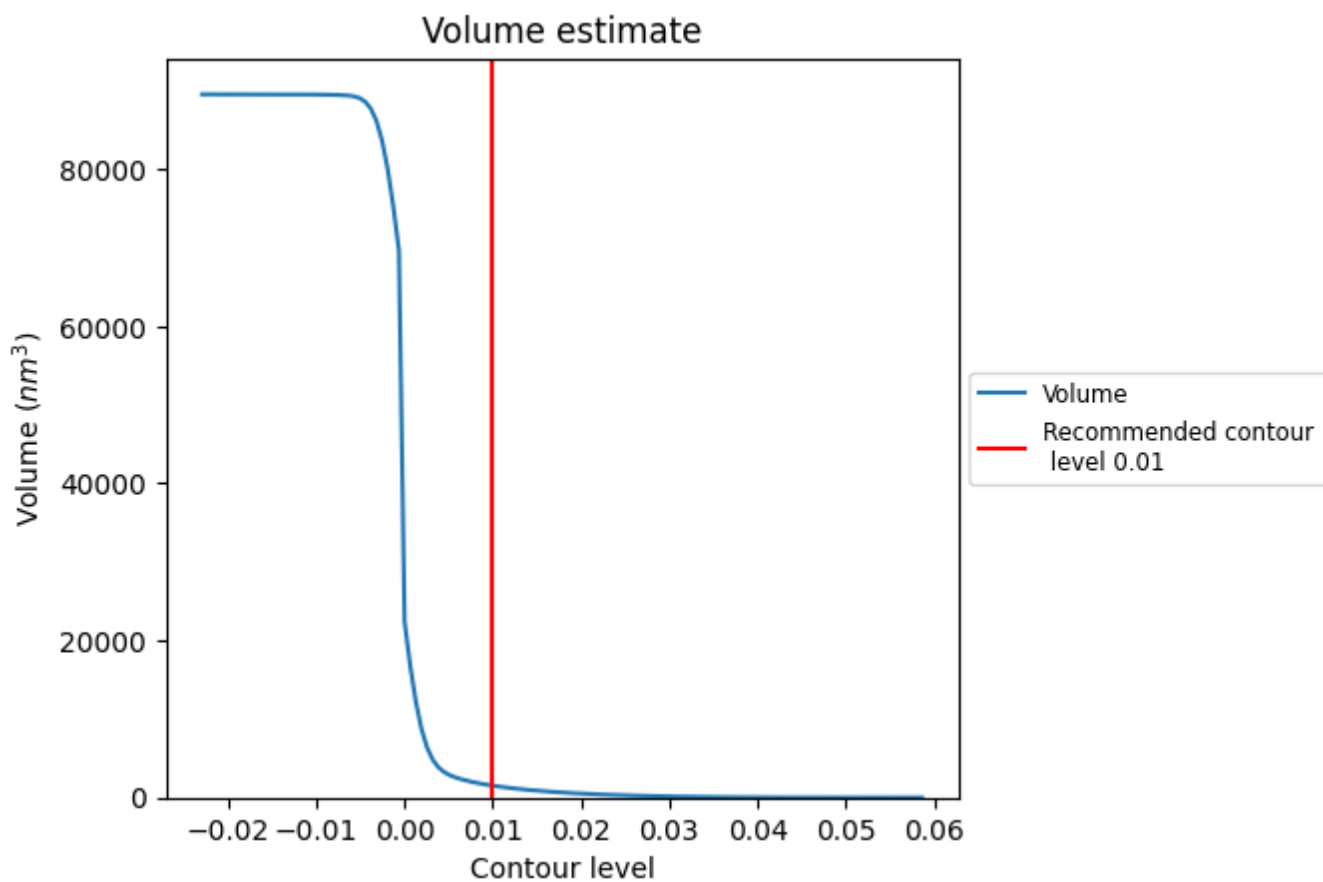
This section contains the results of statistical analysis of the map.

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.

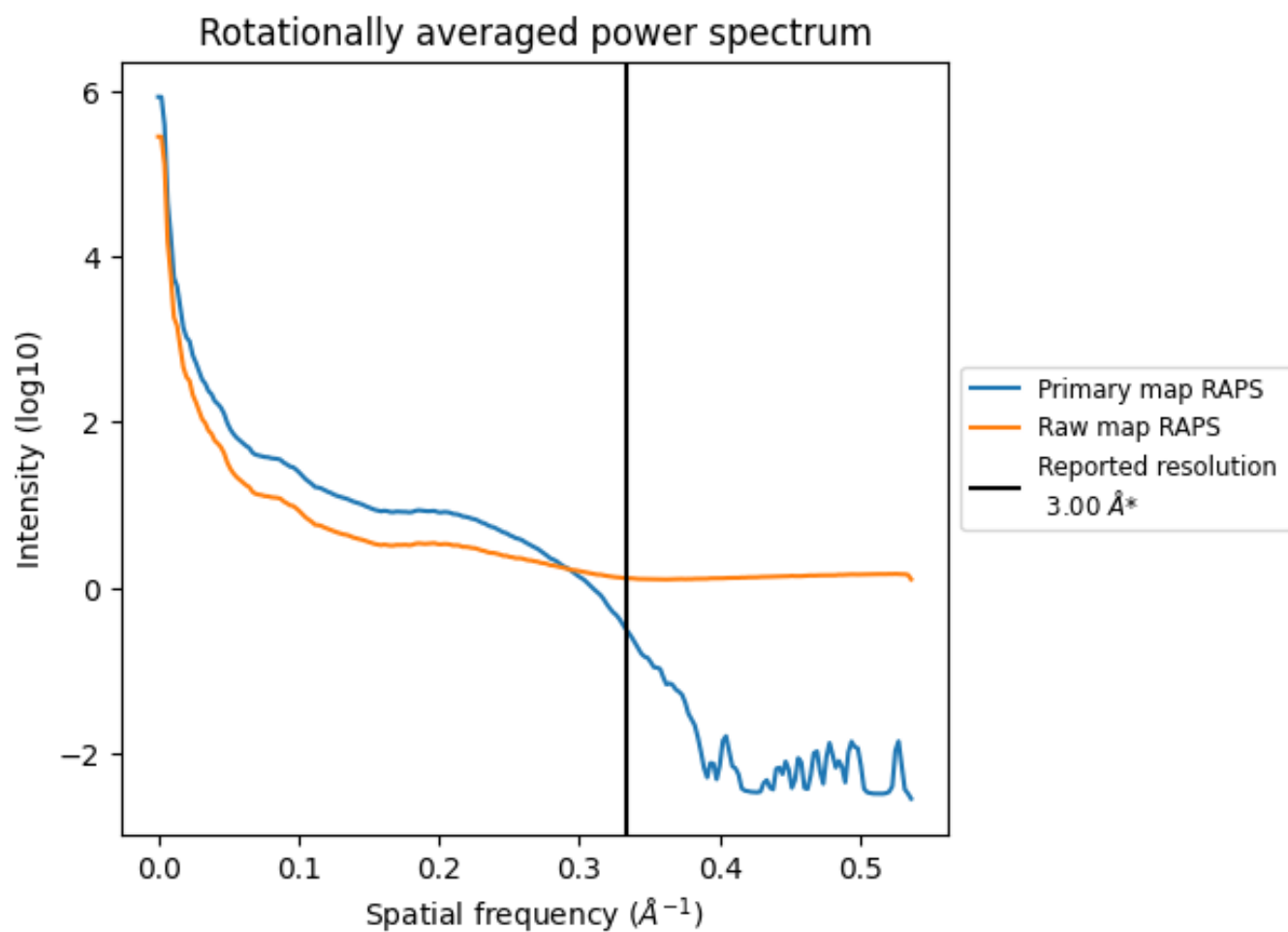
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1517 nm^3 ; this corresponds to an approximate mass of 1370 kDa.

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

7.3 Rotationally averaged power spectrum i

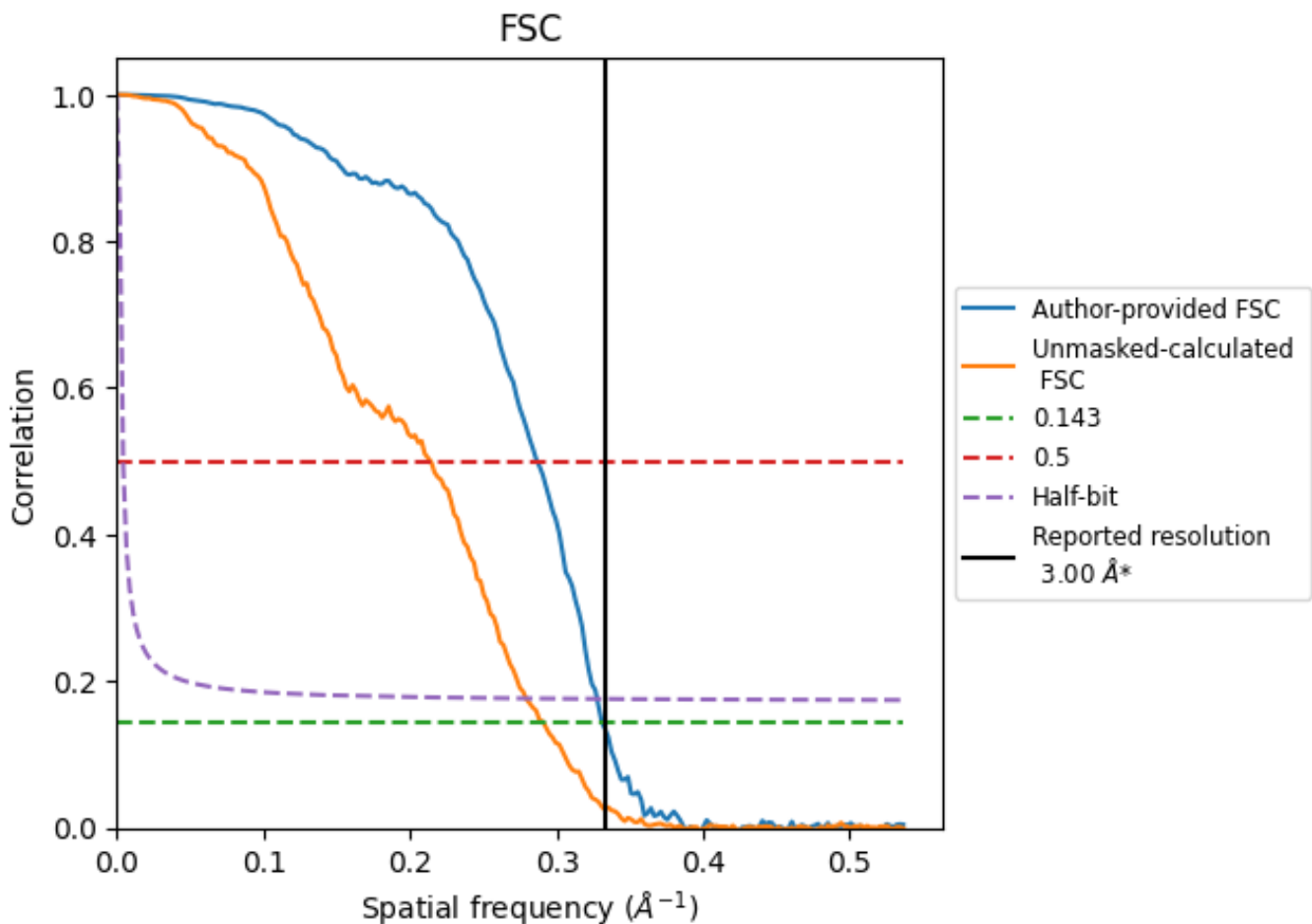


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

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 0.333 Å⁻¹

8.2 Resolution estimates [i](#)

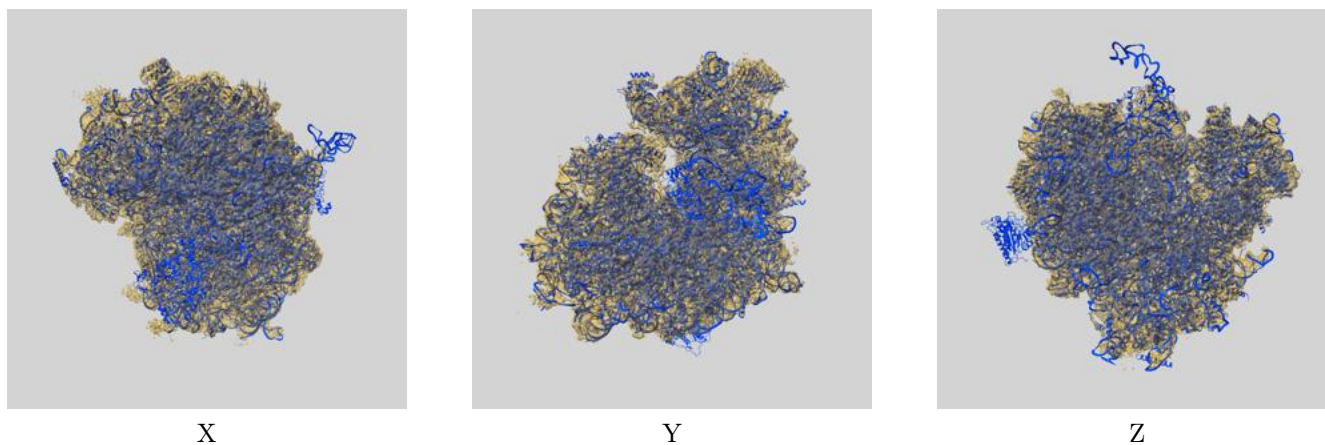
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	3.02	3.49	3.05
Unmasked-calculated*	3.44	4.68	3.58

*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 3.44 differs from the reported value 3.0 by more than 10 %

9 Map-model fit [i](#)

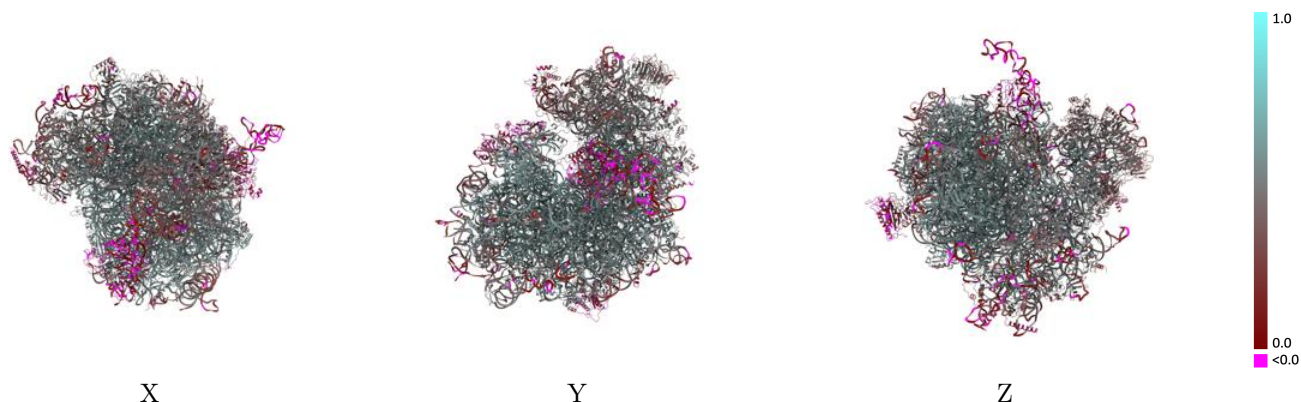
This section contains information regarding the fit between EMDB map EMD-38630 and PDB model 8XSY. Per-residue inclusion information can be found in section 3 on page 22.

9.1 Map-model overlay [i](#)



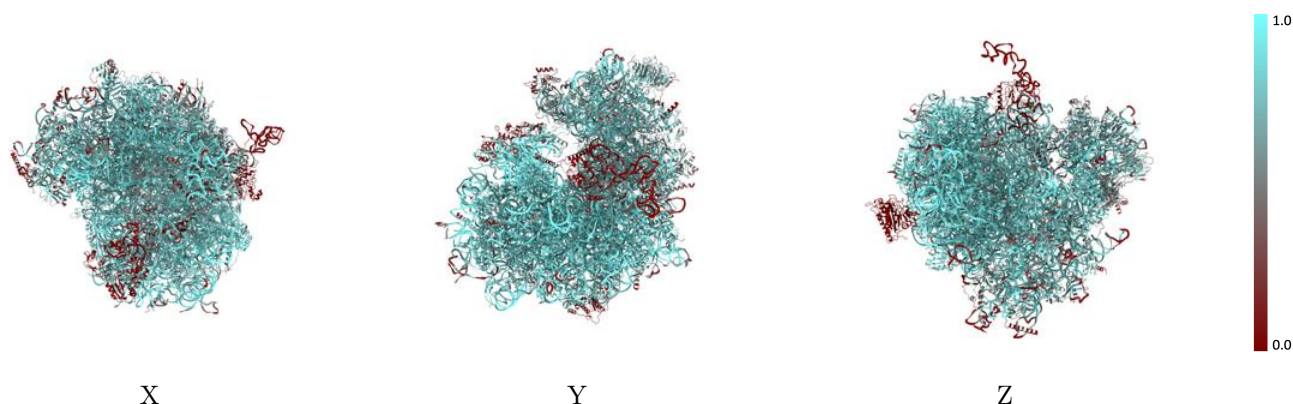
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



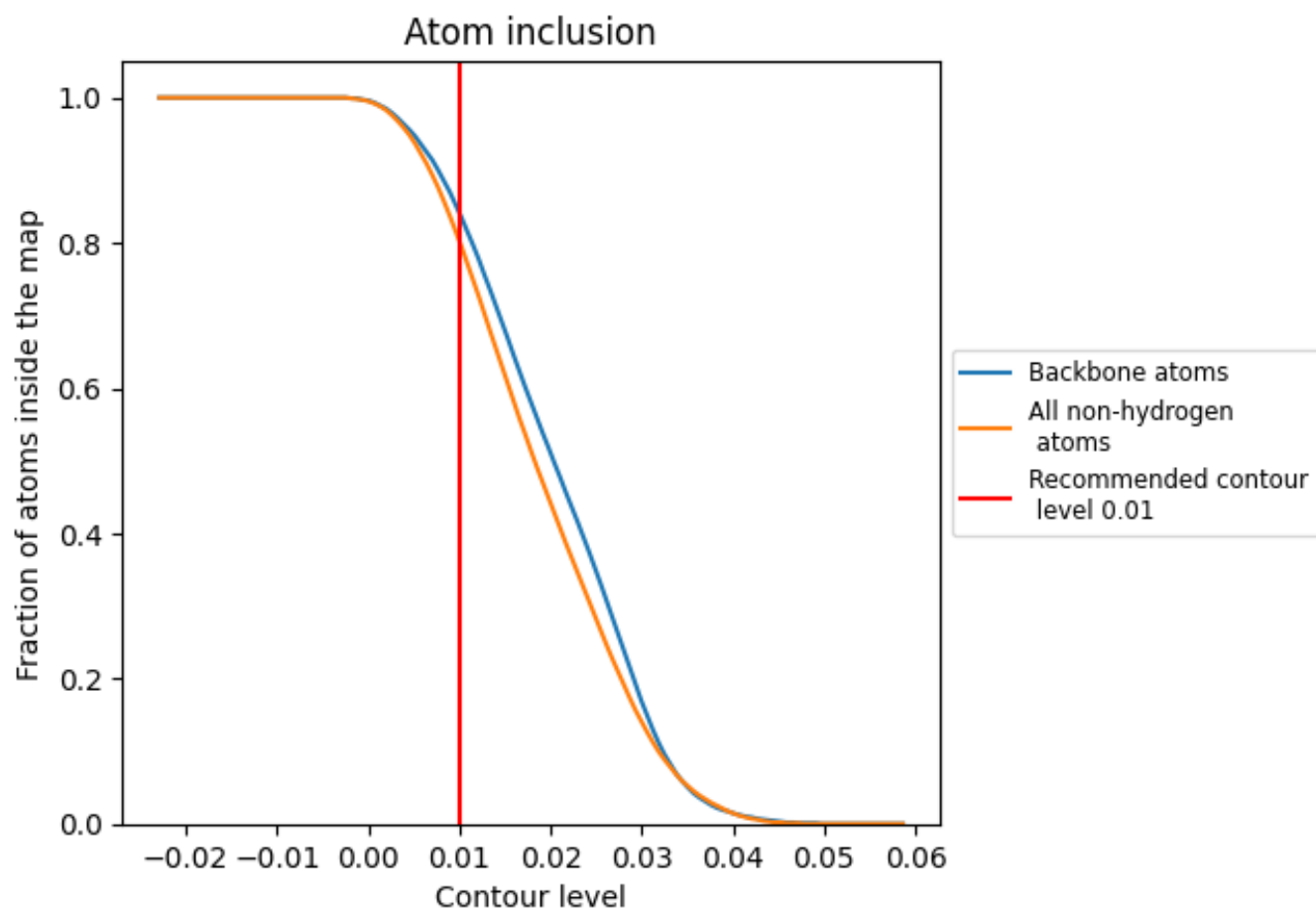
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 80% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8020	 0.4710
CA	 0.0020	 0.1350
CB	 0.2320	 0.4220
CC	 0.2560	 0.1910
CE	 0.0450	 0.3010
L5	 0.8910	 0.4920
L7	 0.9830	 0.5720
L8	 0.9250	 0.5100
LA	 0.8840	 0.5740
LB	 0.8630	 0.5570
LC	 0.8330	 0.5320
LD	 0.8650	 0.5340
LE	 0.8300	 0.5210
LF	 0.8570	 0.5530
LG	 0.7530	 0.4850
LH	 0.8430	 0.5380
LI	 0.8690	 0.5600
LJ	 0.7400	 0.4710
LL	 0.8160	 0.5110
LM	 0.8800	 0.5450
LN	 0.8850	 0.5660
LO	 0.8690	 0.5570
LP	 0.8760	 0.5620
LQ	 0.8620	 0.5600
LR	 0.7930	 0.5050
LS	 0.9060	 0.5760
LT	 0.8460	 0.5440
LU	 0.7840	 0.4790
LV	 0.8590	 0.5560
LW	 0.5080	 0.3340
LX	 0.8180	 0.5340
LY	 0.8450	 0.5300
LZ	 0.8430	 0.5380
La	 0.8790	 0.5660
Lb	 0.7670	 0.4820























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Chain	Atom inclusion	Q-score
Lc	0.8340	0.5300
Ld	0.8330	0.5370
Le	0.8500	0.5600
Lf	0.9030	0.5750
Lg	0.8240	0.5380
Lh	0.8190	0.5230
Li	0.8260	0.5140
Lj	0.8840	0.5560
Lk	0.7810	0.4960
Ll	0.8530	0.5400
Lm	0.8470	0.5400
Ln	0.8570	0.5460
Lo	0.8240	0.5490
Lp	0.8190	0.5520
Lr	0.8430	0.5390
Ls	0.0920	0.0590
Lt	0.0290	0.0220
Lz	0.0250	0.0820
S2	0.8590	0.4560
SA	0.7510	0.4660
SB	0.6940	0.4540
SC	0.7480	0.4690
SD	0.5780	0.3950
SE	0.6760	0.4490
SF	0.6230	0.4060
SG	0.6020	0.3560
SH	0.6260	0.4000
SI	0.6710	0.4530
SJ	0.6780	0.4330
SK	0.5690	0.3300
SL	0.7210	0.4930
SM	0.2910	0.1970
SN	0.7570	0.5010
SO	0.7220	0.4720
SP	0.5560	0.3730
SQ	0.6680	0.4370
SR	0.6670	0.4300
SS	0.5820	0.3730
ST	0.6500	0.4020
SU	0.6000	0.3970
SV	0.7400	0.4710
SW	0.7450	0.4910

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Chain	Atom inclusion	Q-score
SX	 0.7660	 0.4820
SY	 0.6270	 0.3870
SZ	 0.4890	 0.3190
Sa	 0.7770	 0.4940
Sb	 0.6890	 0.4470
Sc	 0.5490	 0.4230
Sd	 0.7280	 0.4510
Se	 0.5590	 0.3840
Sf	 0.3130	 0.2120
Sg	 0.6120	 0.3780