



## Full wwPDB EM Validation Report ⓘ

Dec 17, 2022 – 09:44 pm GMT

PDB ID : 6ZM7  
EMDB ID : EMD-11288  
Title : SARS-CoV-2 Nsp1 bound to the human CCDC124-80S-EBP1 ribosome complex  
Authors : Thoms, M.; Buschauer, R.; Ameismeier, M.; Denk, T.; Kratzat, H.; Mackens-Kiani, T.; Cheng, J.; Berninghausen, O.; Becker, T.; Beckmann, R.  
Deposited on : 2020-07-01  
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

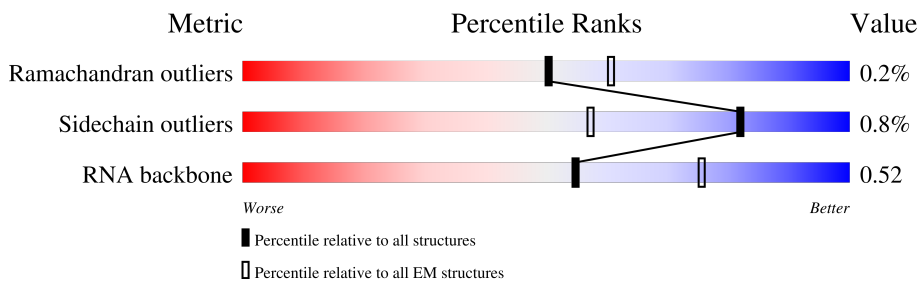
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.3

# 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 2.70 Å.

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



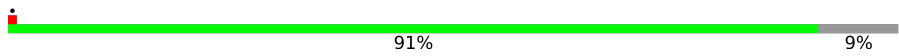

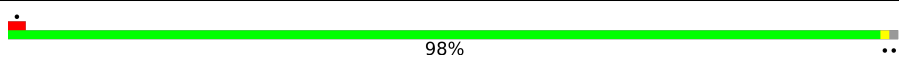
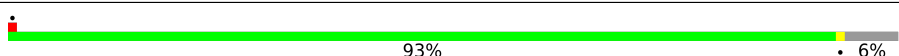
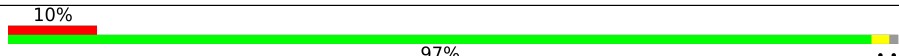
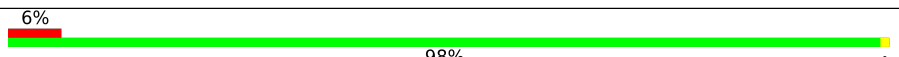
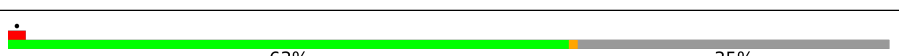
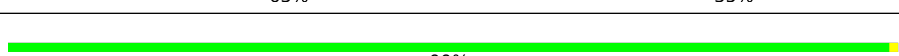
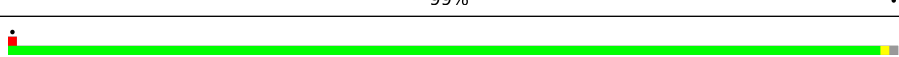
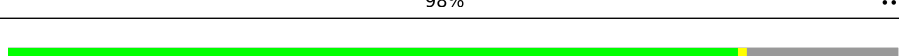
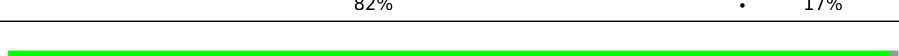
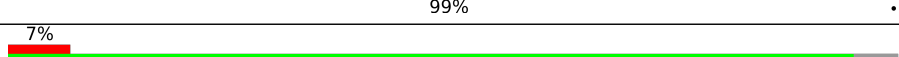
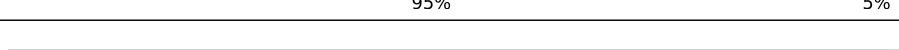
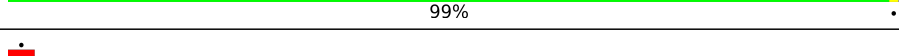
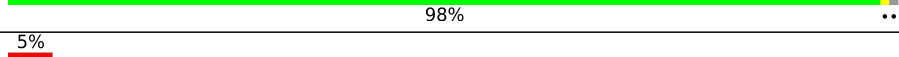

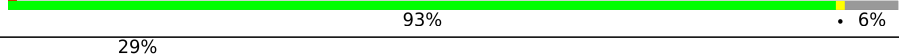


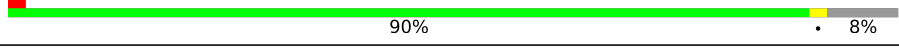
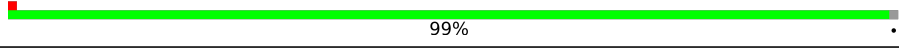
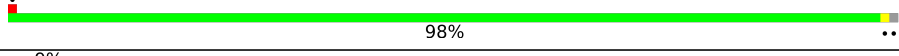



Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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

Mol	Chain	Length	Quality of chain
1	L5	5066	
2	L7	121	
3	L8	157	
4	LA	257	
5	LB	403	
6	LC	427	
7	LD	297	
8	LE	288	

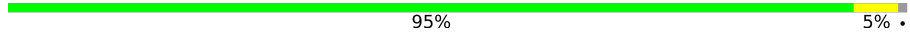
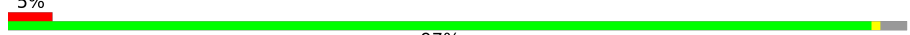
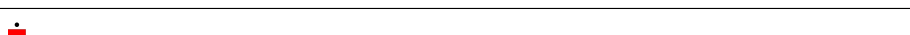
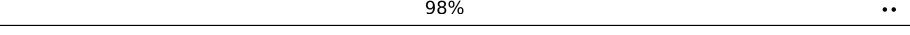
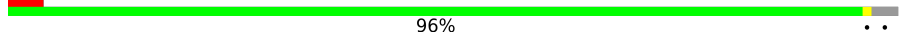

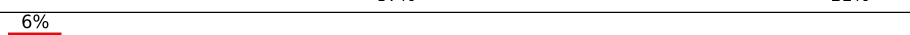
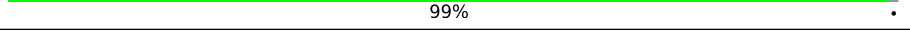
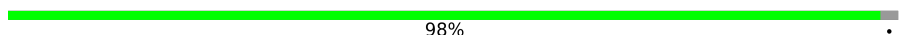

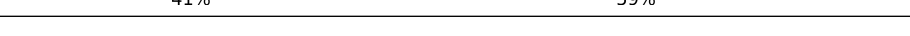
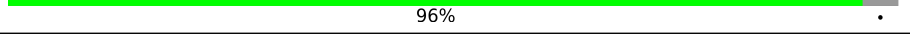
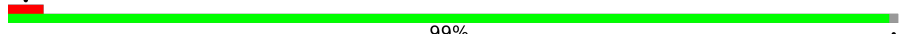
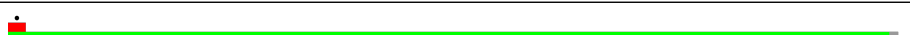
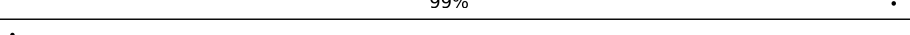



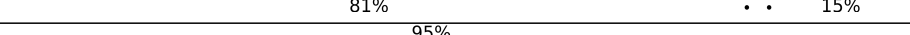


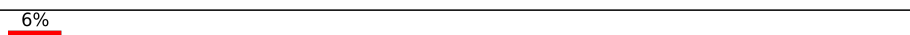



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Mol	Chain	Length	Quality of chain
9	LF	248	
10	LG	266	
11	LH	192	
12	LI	214	
13	LJ	178	
14	LL	211	
15	LM	215	
16	LN	204	
17	LO	203	
18	LP	184	
19	LQ	188	
20	LR	196	
21	LS	176	
22	LT	160	
23	LU	128	
24	LV	140	
25	LW	157	
26	LX	156	
27	LY	145	
28	LZ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

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Mol	Chain	Length	Quality of chain
34	Lf	110	 95% 5%
35	Lg	117	 5% 97%
36	Lh	123	 98%
37	Li	105	 96%
38	Lj	97	 87% 11%
39	Lk	70	 6% 99%
40	Ll	51	 98%
41	Lm	128	 41% 59%
42	Ln	25	 96%
43	Lo	106	 99%
44	Lp	92	 99%
45	Lr	137	 91% 9%
46	Ls	317	 60% 60% 38%
47	Lt	165	 85% 81% 15%
48	Lz	217	 95% 95% 5%
49	S2	1869	 7% 69% 22% 7%
50	SA	295	 6% 74% 25%
51	SB	264	 80% 19%
52	SD	243	 10% 91% 7%
53	SE	263	 99%
54	SF	204	 89% 10%
55	SH	194	 20% 94%
56	SI	208	 7% 98%
57	SK	165	 7% 58% 41%
58	SL	158	 11% 96%

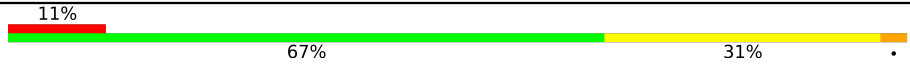


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Mol	Chain	Length	Quality of chain
59	SP	145	15% 88% 11%
60	SQ	146	5% 97%
61	SR	135	10% 99%
62	SS	152	95% 5%
63	ST	145	6% 96%
64	SU	119	17% 87% 13%
65	SV	83	96%
66	SX	143	94%
67	Sa	115	88% 11%
68	Sc	69	12% 91% 7%
69	Sd	56	95%
70	Sg	317	25% 97%
71	SC	293	75% 24%
72	SG	249	15% 93% 5%
73	SJ	194	94% 5%
74	SM	132	78% 89% 8%
75	SN	151	99%
76	SO	151	7% 92% 7%
77	SW	130	99%
78	SY	133	8% 95%
79	SZ	125	8% 55% 40%
80	Sb	84	10% 98%
81	Se	59	19% 97%
82	Sf	156	29% 42% 57%
83	CA	394	88% 89% 10%

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Mol	Chain	Length	Quality of chain
84	CC	75	
85	CE	223	
86	CF	180	

## 2 Entry composition

There are 88 unique types of molecules in this entry. The entry contains 225534 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 ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	L5	3773	80138	35655	14589	26122	3772	0	0

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

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 ribosomal RNA.

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	236	1904	1222	361	317	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 60S ribosomal protein L10-like.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	LI	202	1634	1037	314	269	14	0	0

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

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

- 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
20	LR	187	Total	C	N	O	S	0	0
			1566	971	336	250	9		

- 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	101	825	529	144	150	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	134	1115	700	226	186	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	876	546	189	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
			Total	C	N	O	S		
41	Lm	52	429	266	90	67	6	0	0

- 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 60S acidic ribosomal protein P0.

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
48	Lz	217	Total	C	N	O	S	0	0
			1741	1113	312	307	9		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
49	S2	1740	36899	16459	6598	12103	1739	0	0

- 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
56	SI	206	Total	C	N	O	S	0	0
			1686	1058	332	291	5		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
57	SK	98	Total	C	N	O	S	0	0
			827	539	148	134	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
58	SL	153	Total	C	N	O	S	0	0
			1247	793	234	214	6		

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
62	SS	145	Total	C	N	O	S	0	0
			1198	751	242	203	2		

- 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
			Total	C	N	O	S		
68	Sc	64	506	308	102	94	2	0	0

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

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

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



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

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
74	SM	122	940	590	164	177	9	0	0

- 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	140	1049	642	204	197	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	131	1065	673	209	178	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
			Total	C	N	O	S		
81	Se	58	439	268	97	73	1	0	0

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

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

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

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

- Molecule 84 is a RNA chain called tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
84	CC	75	1589	710	279	525	75	0	0

- Molecule 85 is a protein called Coiled-coil domain-containing protein 124.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
85	CE	124	1043	637	208	194	4	0	0

- Molecule 86 is a protein called Non-structural protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
86	CF	29	239	145	43	50	1	0	0

- 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	
87	LA	1	Total	Mg	0
			1	1	
87	LI	1	Total	Mg	0
			1	1	
87	LP	1	Total	Mg	0
			1	1	
87	LT	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	S2	29	Total	Mg	0
			29	29	
87	SG	1	Total	Mg	0
			1	1	

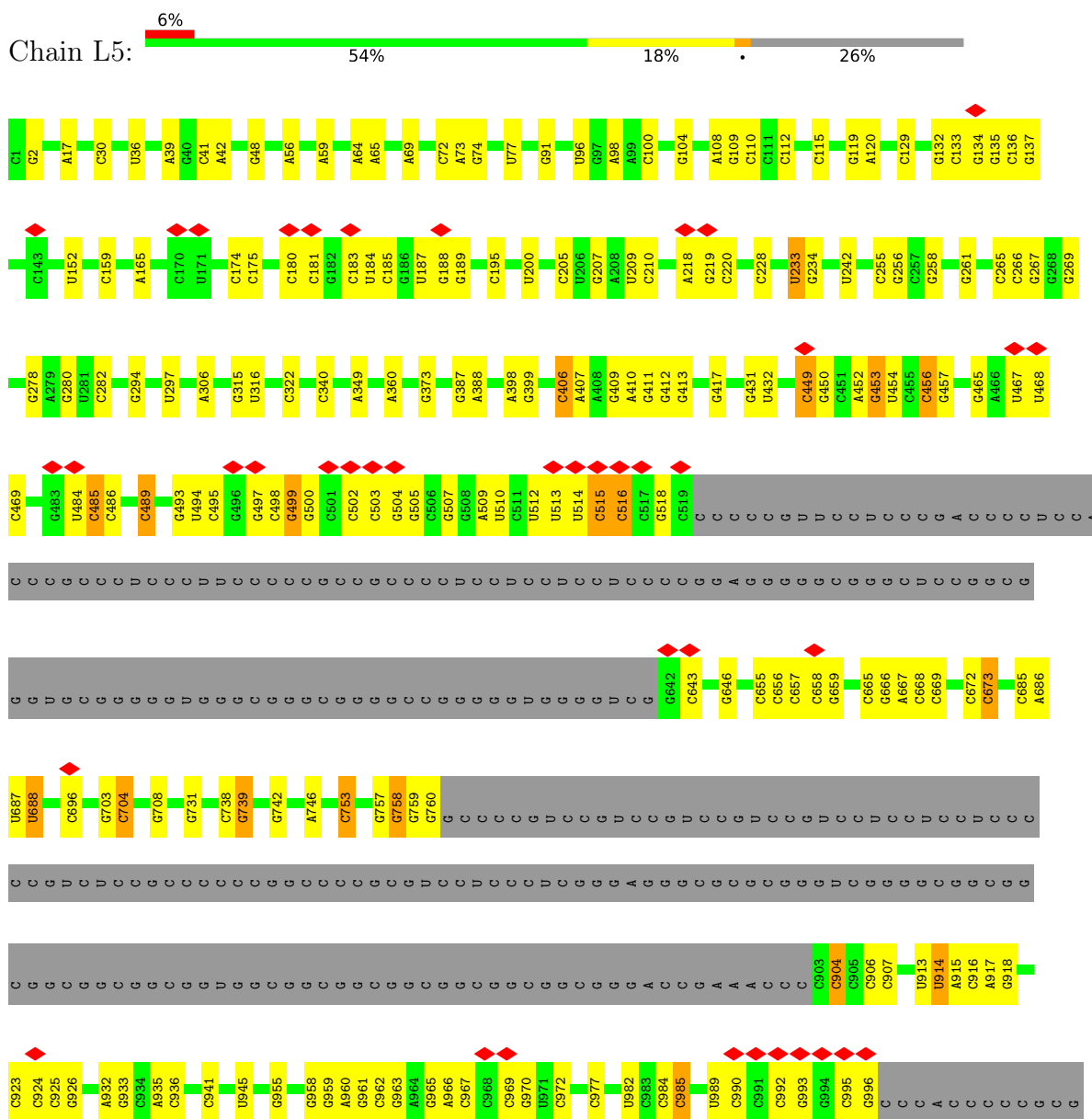
- Molecule 88 is ZINC ION (three-letter code: ZN) (formula: Zn).

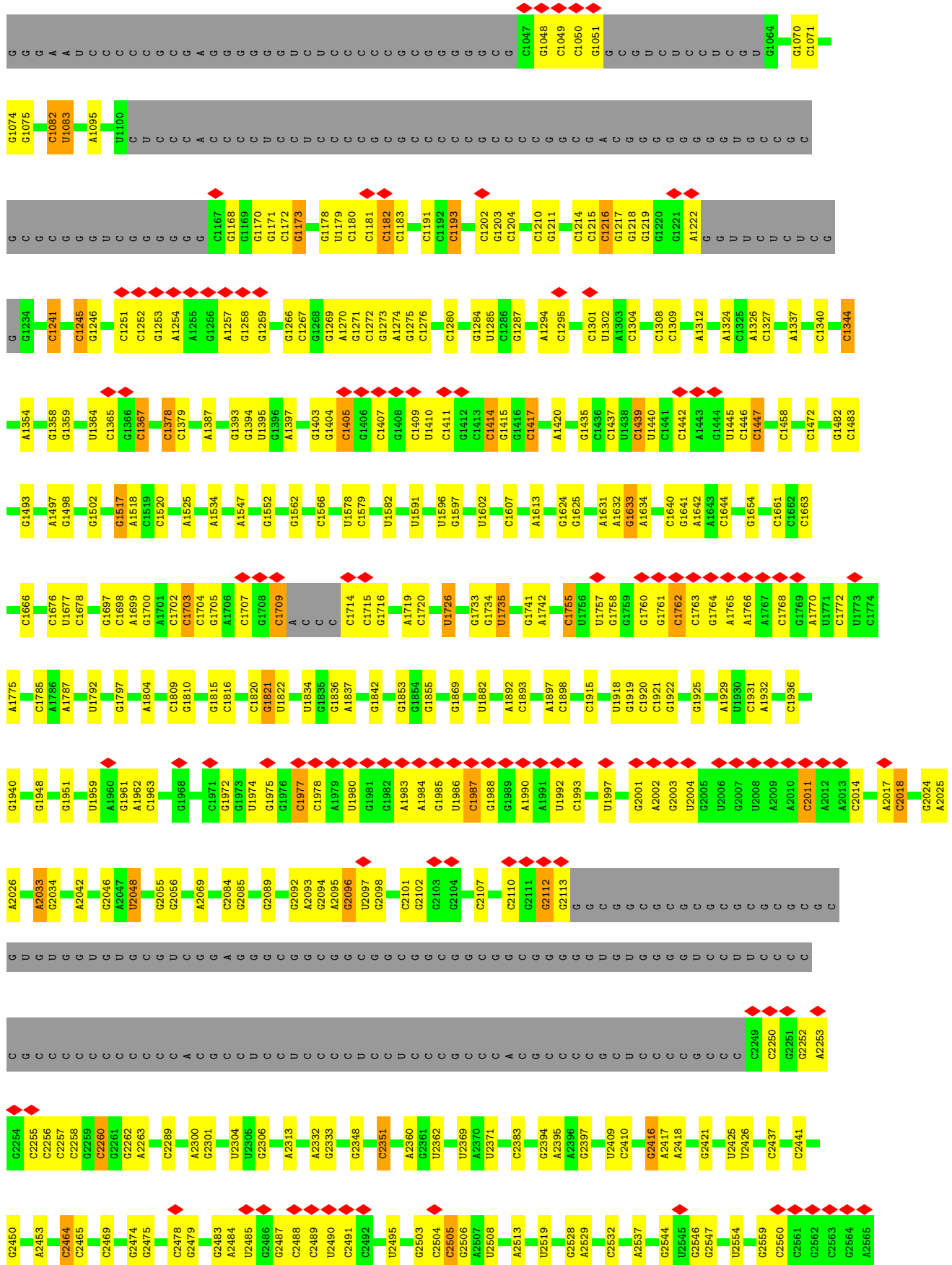
Mol	Chain	Residues	Atoms		AltConf
88	Lg	1	Total 1	Zn 1	0
88	Lj	1	Total 1	Zn 1	0
88	Lm	1	Total 1	Zn 1	0
88	Lo	1	Total 1	Zn 1	0
88	Lp	1	Total 1	Zn 1	0
88	Sa	1	Total 1	Zn 1	0
88	Sd	1	Total 1	Zn 1	0
88	Sf	1	Total 1	Zn 1	0

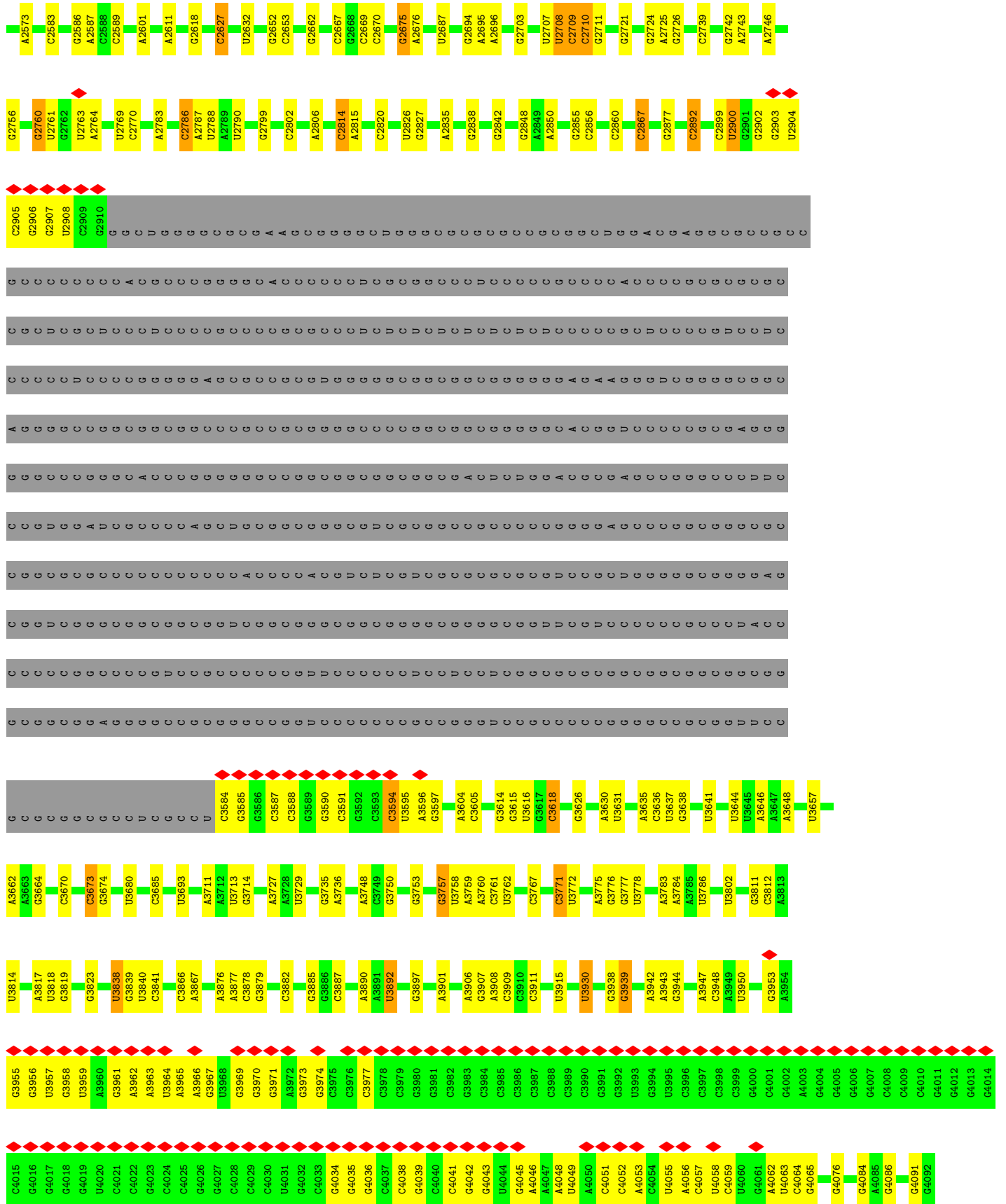
### 3 Residue-property plots [i](#)

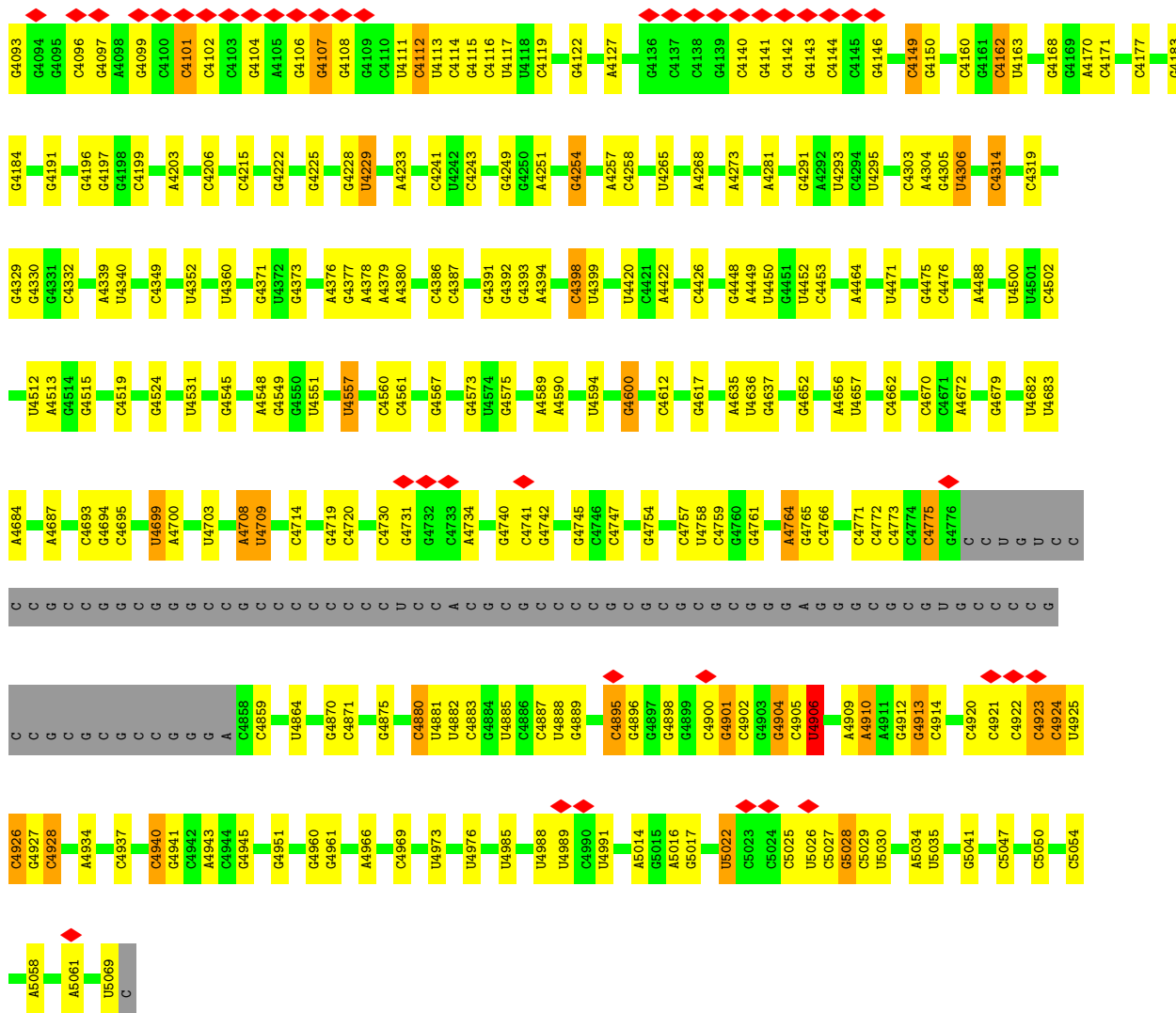
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

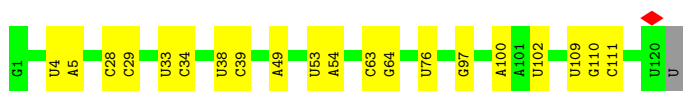
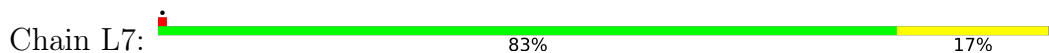




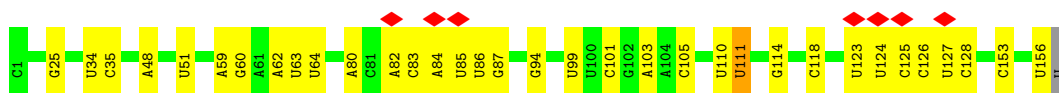
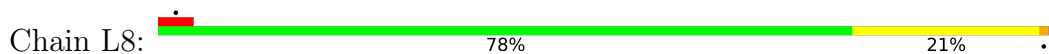




• Molecule 2: 5S ribosomal RNA



• Molecule 3: 5.8S ribosomal RNA



• Molecule 4: 60S ribosomal protein L8



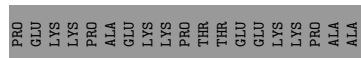
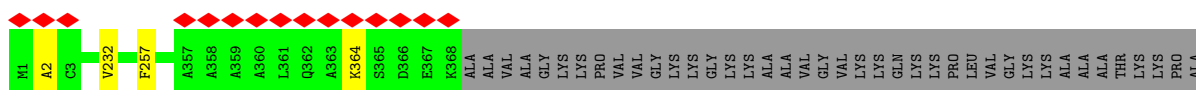
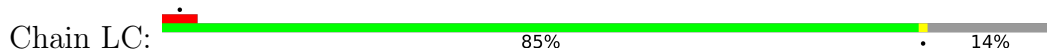




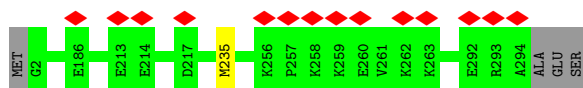
- Molecule 5: 60S ribosomal protein L3



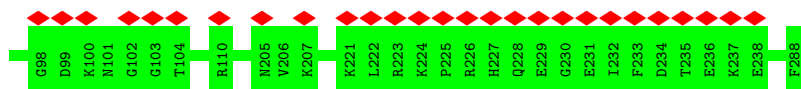
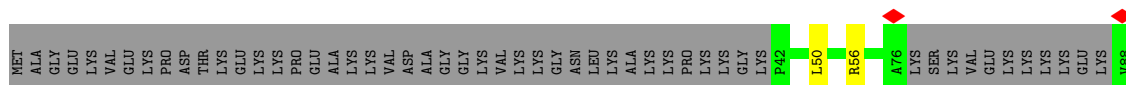
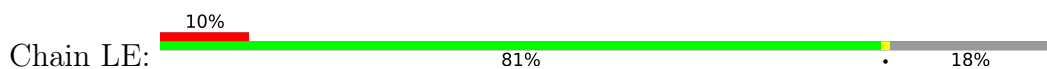
- Molecule 6: 60S ribosomal protein L4



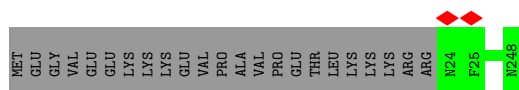
- Molecule 7: 60S ribosomal protein L5



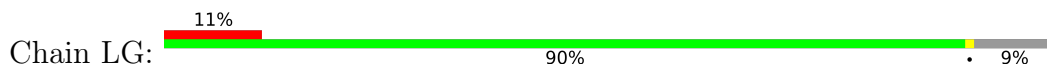
- Molecule 8: 60S ribosomal protein L6

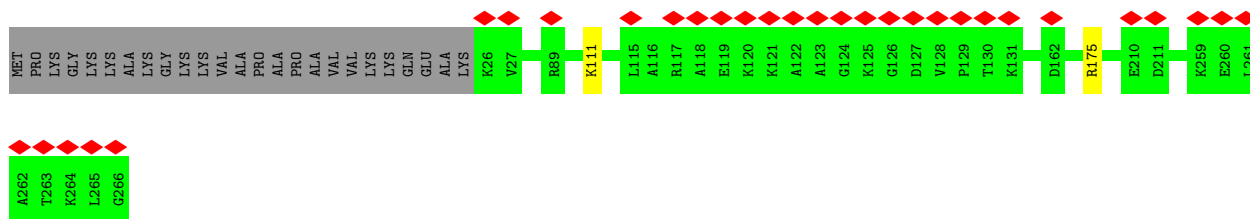


- Molecule 9: 60S ribosomal protein L7

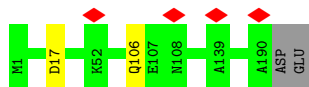


- Molecule 10: 60S ribosomal protein L7a

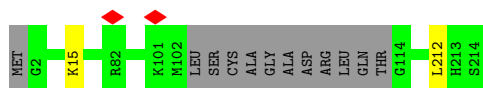




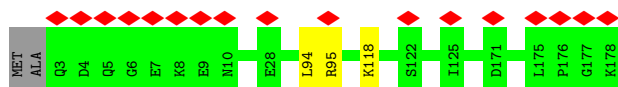
- Molecule 11: 60S ribosomal protein L9



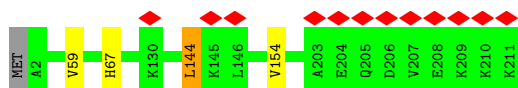
- Molecule 12: 60S ribosomal protein L10-like



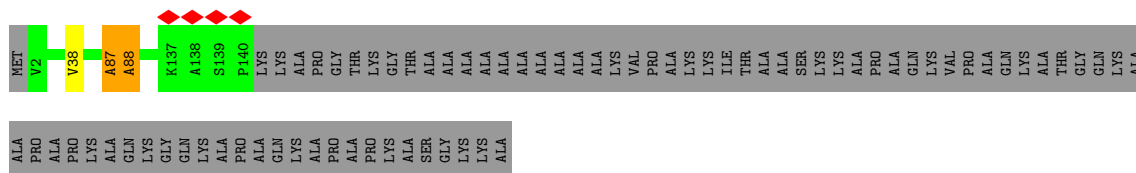
- Molecule 13: 60S ribosomal protein L11



- Molecule 14: 60S ribosomal protein L13



- Molecule 15: 60S ribosomal protein L14



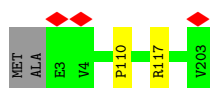
- Molecule 16: 60S ribosomal protein L15

Chain LN:  99%





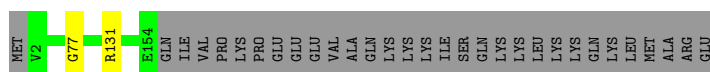
- Molecule 17: 60S ribosomal protein L13a

Chain LO:  98%



- Molecule 18: 60S ribosomal protein L17

Chain LP:  82%  17%



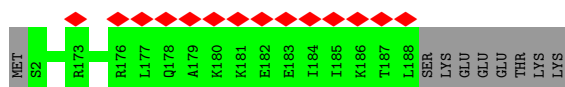
- Molecule 19: 60S ribosomal protein L18

Chain LQ:  99%



- Molecule 20: 60S ribosomal protein L19

Chain LR:  7%  95%  5%



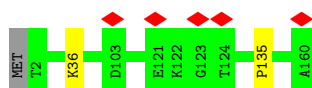
- Molecule 21: 60S ribosomal protein L18a

Chain LS:  99%

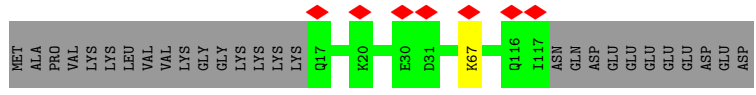
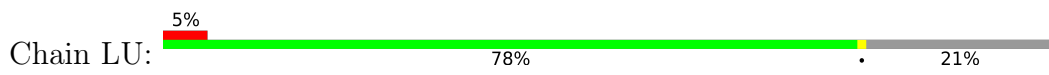


- Molecule 22: 60S ribosomal protein L21

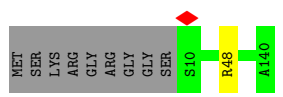
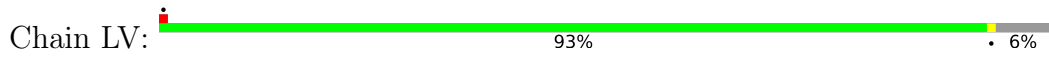
Chain LT:  98%



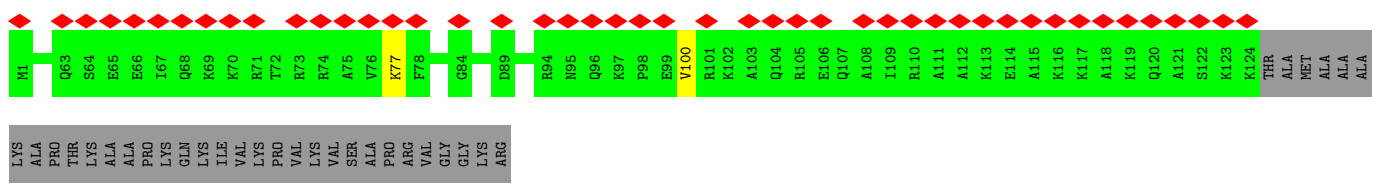
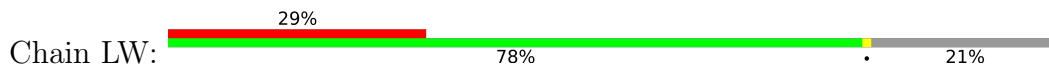
- Molecule 23: 60S ribosomal protein L22



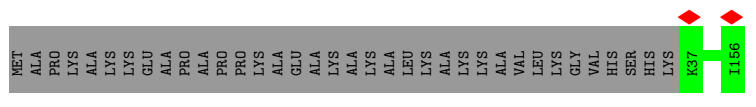
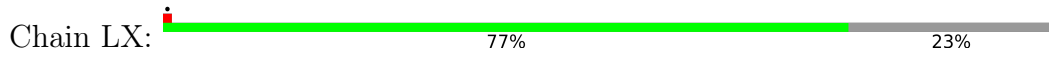
• Molecule 24: 60S ribosomal protein L23



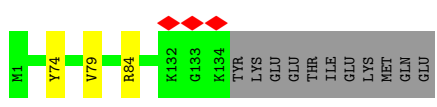
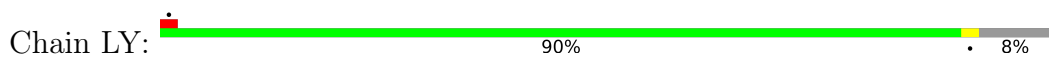
• Molecule 25: 60S ribosomal protein L24



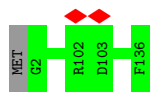
• Molecule 26: 60S ribosomal protein L23a



• Molecule 27: 60S ribosomal protein L26

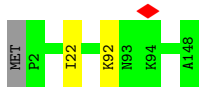


• Molecule 28: 60S ribosomal protein L27

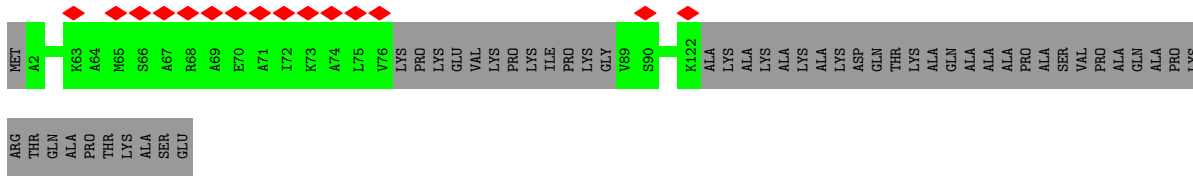
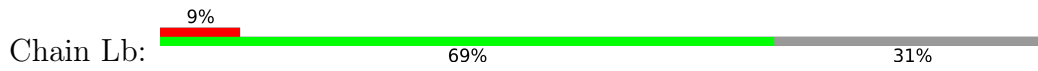


• Molecule 29: 60S ribosomal protein L27a

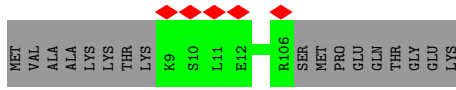
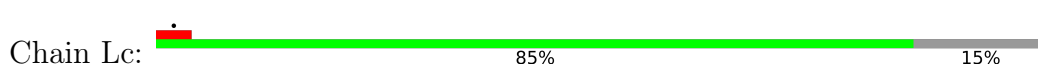




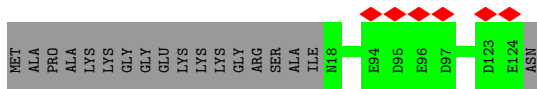
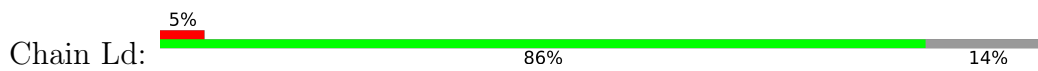
- Molecule 30: 60S ribosomal protein L29



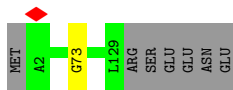
- Molecule 31: 60S ribosomal protein L30



- Molecule 32: 60S ribosomal protein L31



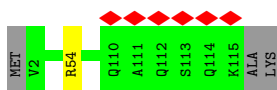
- Molecule 33: 60S ribosomal protein L32



- Molecule 34: 60S ribosomal protein L35a

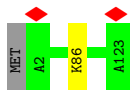


- Molecule 35: 60S ribosomal protein L34



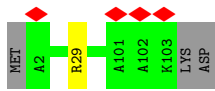
- Molecule 36: 60S ribosomal protein L35

Chain Lh:  98%




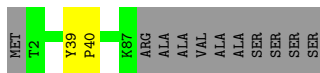
- Molecule 37: 60S ribosomal protein L36

Chain Li:  96%



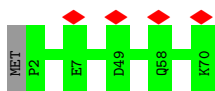
- Molecule 38: 60S ribosomal protein L37

Chain Lj:  87% 11%



- Molecule 39: 60S ribosomal protein L38

Chain Lk:  6% 99%



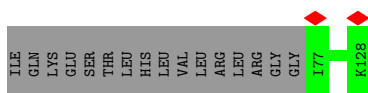
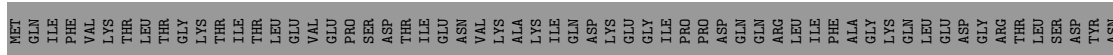
- Molecule 40: 60S ribosomal protein L39

Chain Ll:  98%



- Molecule 41: Ubiquitin-60S ribosomal protein L40

Chain Lm:  41% 59%



- Molecule 42: 60S ribosomal protein L41

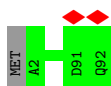
Chain Ln:  96%



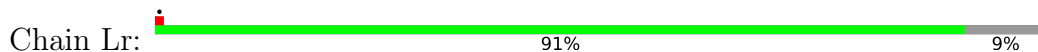
- Molecule 43: 60S ribosomal protein L36a



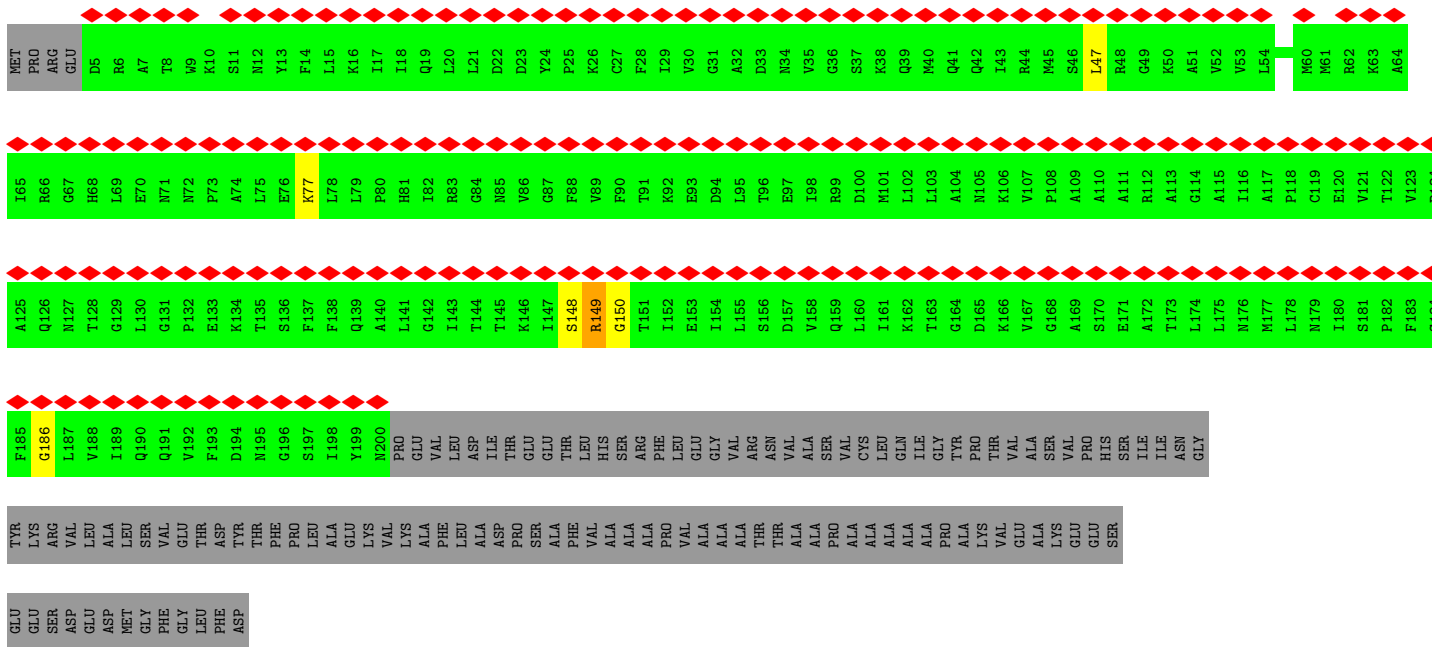
- Molecule 44: 60S ribosomal protein L37a



- Molecule 45: 60S ribosomal protein L28



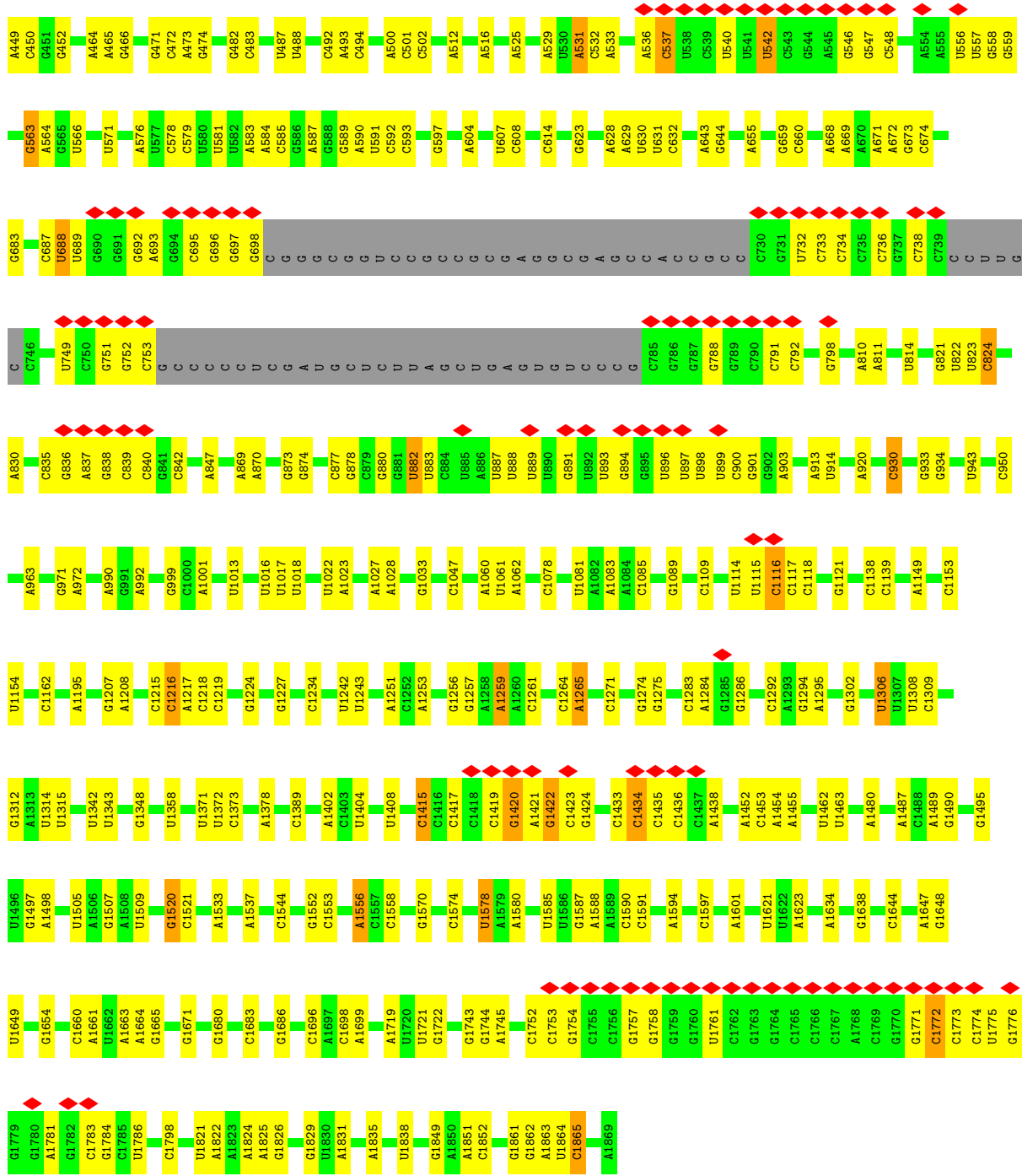
- Molecule 46: 60S acidic ribosomal protein P0



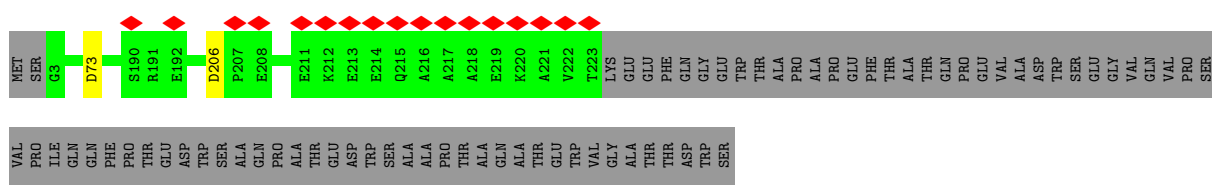
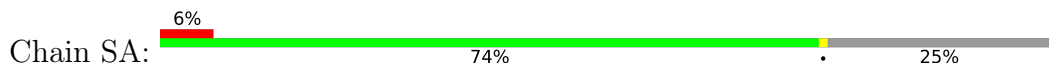
- Molecule 47: 60S ribosomal protein L12



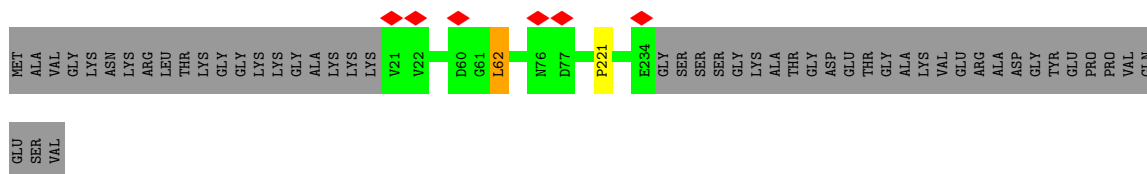
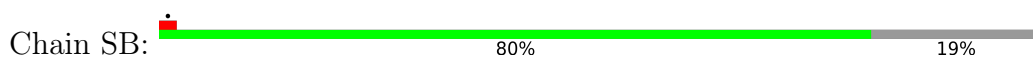




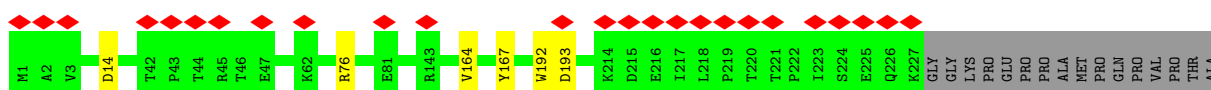
• Molecule 50: 40S ribosomal protein SA



• Molecule 51: 40S ribosomal protein S3a



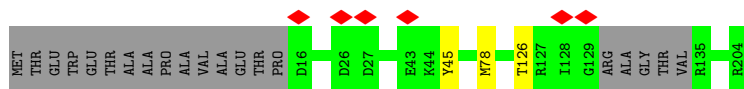
• Molecule 52: 40S ribosomal protein S3



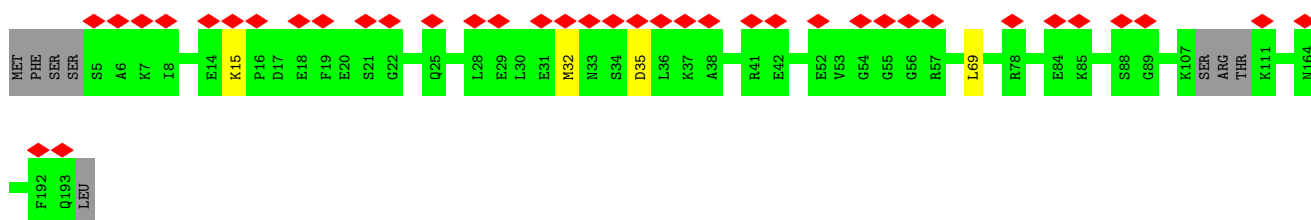
• Molecule 53: 40S ribosomal protein S4, X isoform



• Molecule 54: 40S ribosomal protein S5

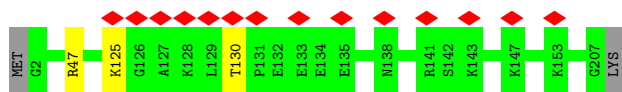


• Molecule 55: 40S ribosomal protein S7

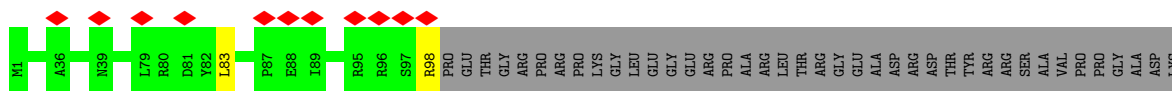


• Molecule 56: 40S ribosomal protein S8

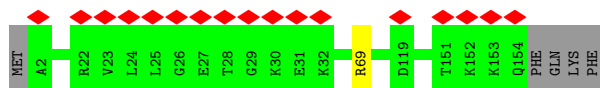




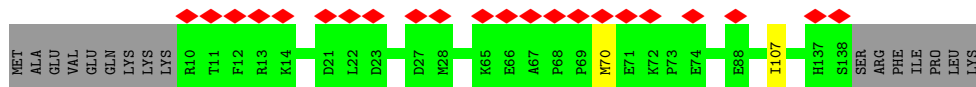
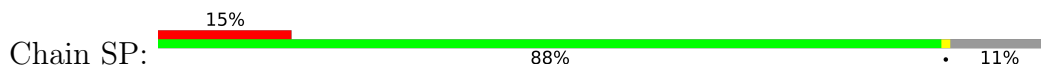
- Molecule 57: 40S ribosomal protein S10



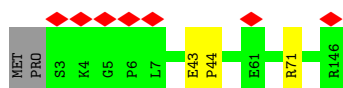
- Molecule 58: 40S ribosomal protein S11



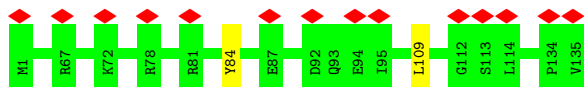
- Molecule 59: 40S ribosomal protein S15



- Molecule 60: 40S ribosomal protein S16

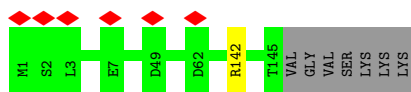


- Molecule 61: 40S ribosomal protein S17

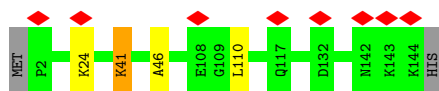


- Molecule 62: 40S ribosomal protein S18

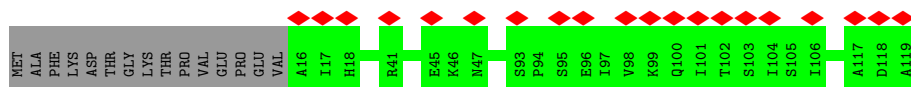
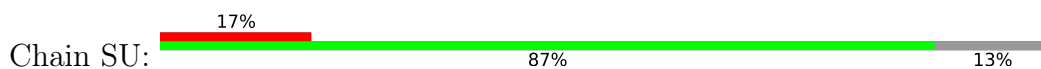




- Molecule 63: 40S ribosomal protein S19



- Molecule 64: 40S ribosomal protein S20



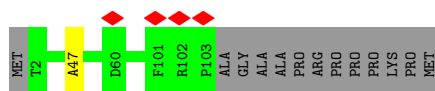
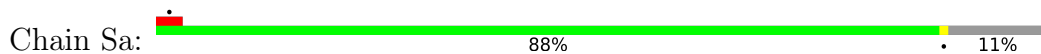
- Molecule 65: 40S ribosomal protein S21



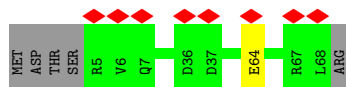
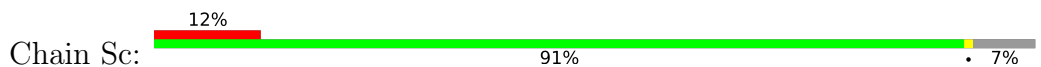
- Molecule 66: 40S ribosomal protein S23



- Molecule 67: 40S ribosomal protein S26



- Molecule 68: 40S ribosomal protein S28



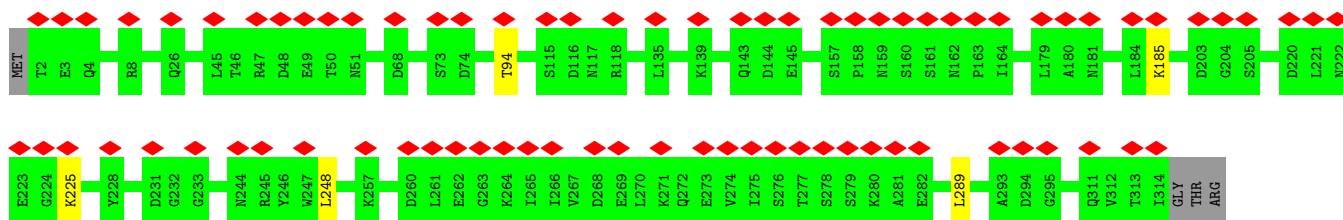
- Molecule 69: 40S ribosomal protein S29

Chain Sd:  95%




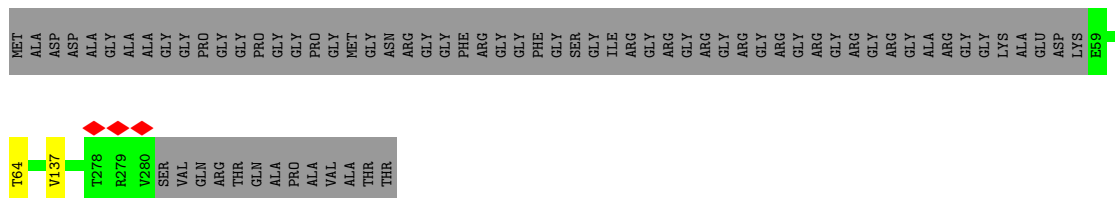
• Molecule 70: Receptor of activated protein C kinase 1

Chain Sg:  97%

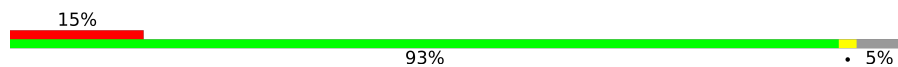


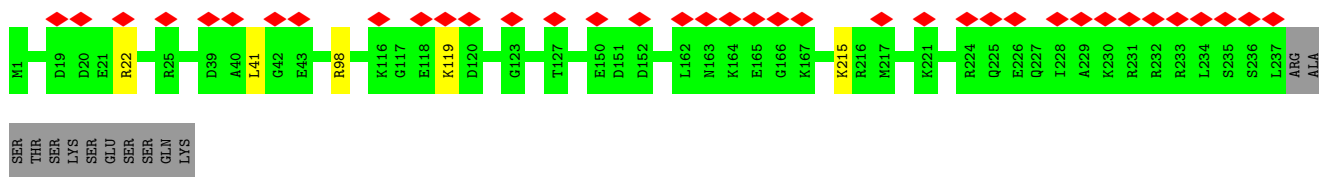
• Molecule 71: 40S ribosomal protein S2

Chain SC:  75%



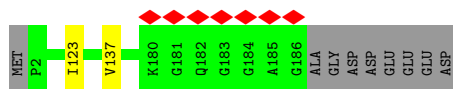
• Molecule 72: 40S ribosomal protein S6

Chain SG:  93%

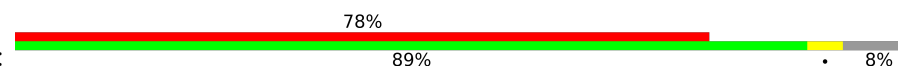


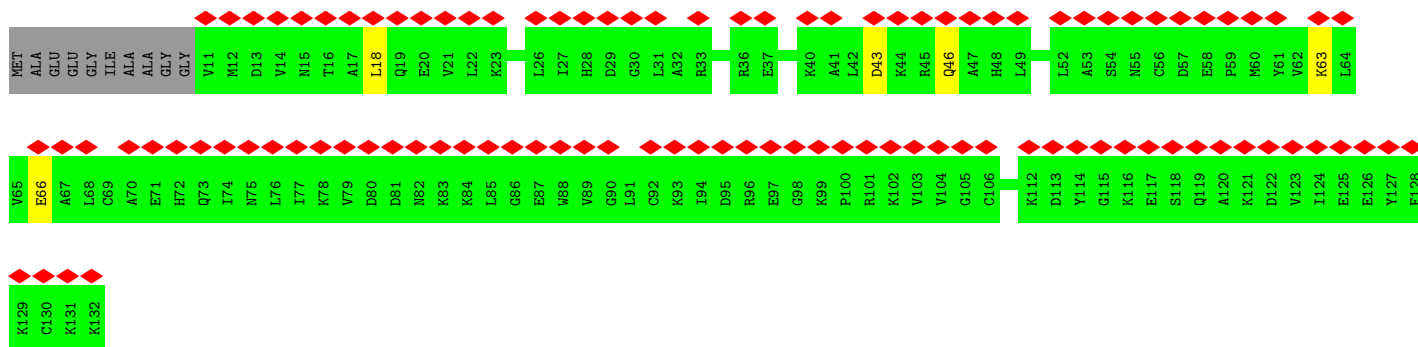
• Molecule 73: 40S ribosomal protein S9

Chain SJ:  94%



• Molecule 74: 40S ribosomal protein S12

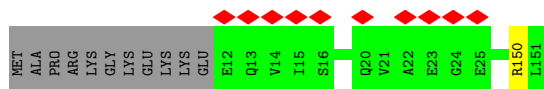
Chain SM:  89%



• Molecule 75: 40S ribosomal protein S13



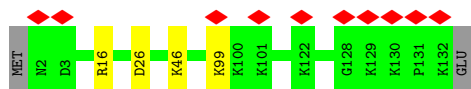
• Molecule 76: 40S ribosomal protein S14



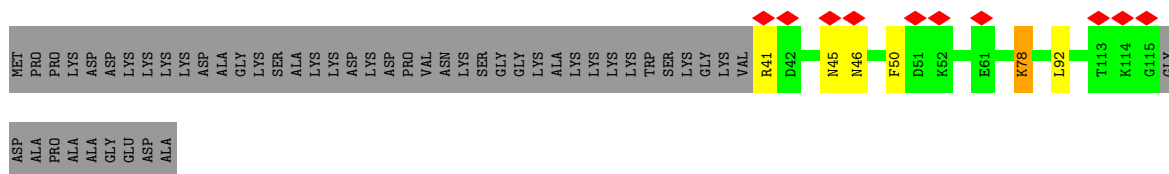
• Molecule 77: 40S ribosomal protein S15a



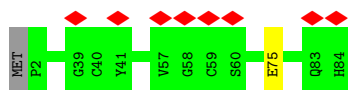
• Molecule 78: 40S ribosomal protein S24



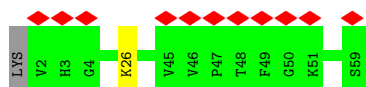
• Molecule 79: 40S ribosomal protein S25



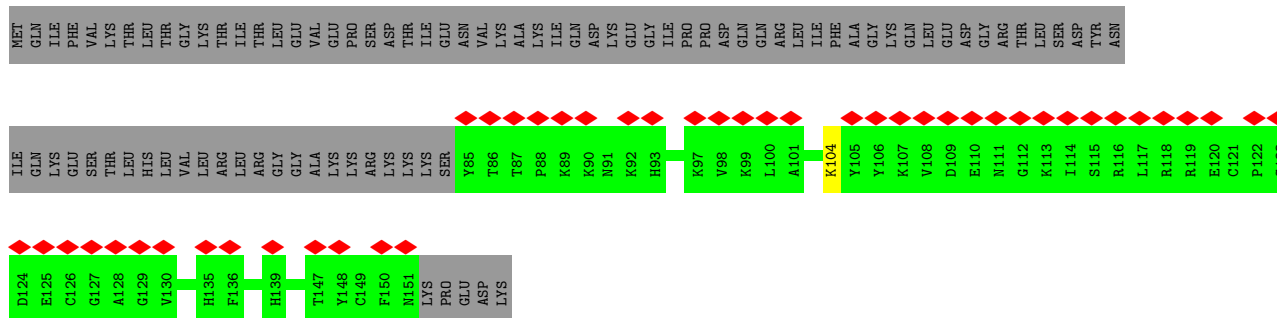
• Molecule 80: 40S ribosomal protein S27



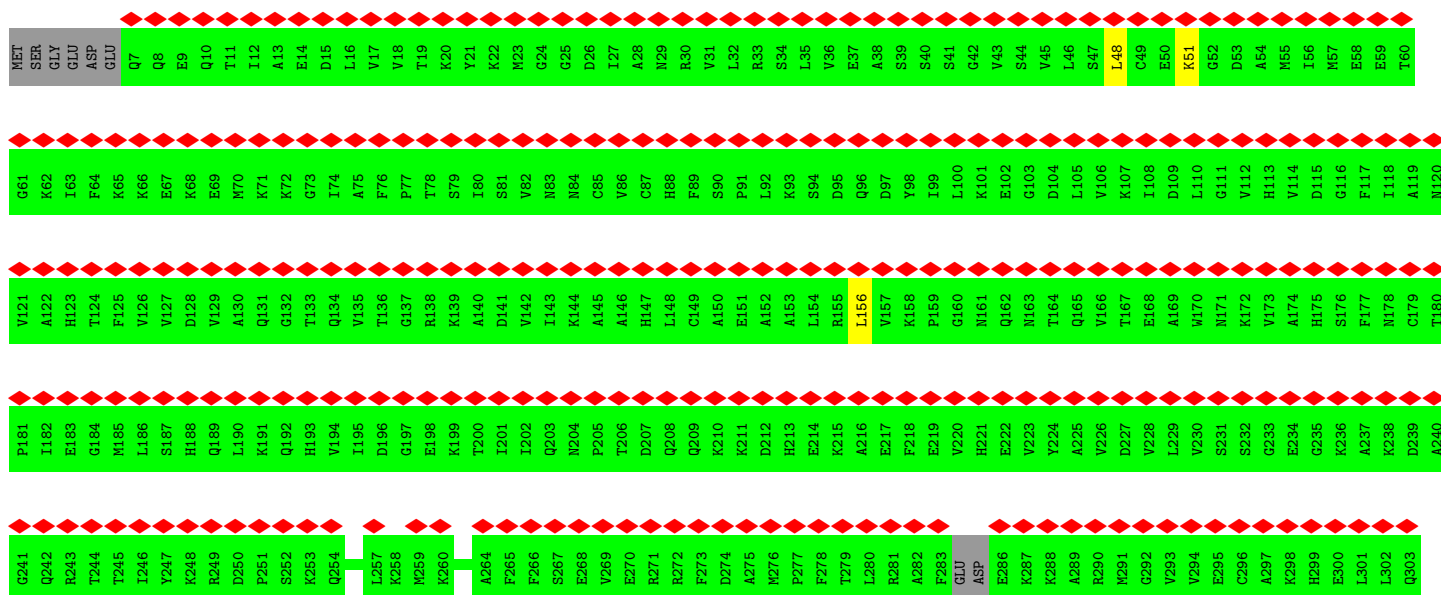
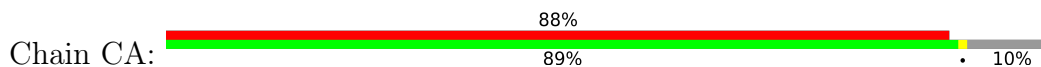
• Molecule 81: 40S ribosomal protein S30



• Molecule 82: Ubiquitin-40S ribosomal protein S27a



• Molecule 83: Proliferation-associated protein 2G4







## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	52512	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	44.8	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.527	Depositor
Minimum map value	-0.124	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.019	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	423.6, 423.6, 423.6	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.059, 1.059, 1.059	Depositor

## 5 Model quality i

### 5.1 Standard geometry i

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	L5	0.77	1/89595 (0.0%)	1.12	556/139686 (0.4%)
2	L7	0.70	0/2861	1.05	12/4459 (0.3%)
3	L8	0.76	0/3701	1.05	15/5766 (0.3%)
4	LA	0.45	0/1936	0.66	1/2596 (0.0%)
5	LB	0.41	0/3306	0.61	3/4424 (0.1%)
6	LC	0.41	0/2981	0.58	1/4002 (0.0%)
7	LD	0.38	0/2428	0.54	1/3252 (0.0%)
8	LE	0.35	0/1942	0.60	1/2606 (0.0%)
9	LF	0.41	0/1905	0.54	0/2539
10	LG	0.38	0/1960	0.54	0/2637
11	LH	0.38	0/1537	0.55	0/2066
12	LI	0.40	0/1673	0.56	1/2233 (0.0%)
13	LJ	0.34	0/1433	0.61	0/1915
14	LL	0.38	0/1732	0.55	1/2315 (0.0%)
15	LM	0.38	0/1161	0.57	1/1554 (0.1%)
16	LN	0.44	0/1746	0.57	0/2338
17	LO	0.40	0/1682	0.50	0/2250
18	LP	0.41	0/1268	0.54	0/1701
19	LQ	0.40	0/1537	0.53	0/2052
20	LR	0.37	0/1582	0.50	0/2091
21	LS	0.41	0/1493	0.52	0/2003
22	LT	0.41	0/1326	0.57	0/1770
23	LU	0.38	0/839	0.66	0/1126
24	LV	0.40	0/993	0.57	0/1332
25	LW	0.36	0/1030	0.53	0/1364
26	LX	0.38	0/1002	0.50	0/1345
27	LY	0.40	0/1132	0.55	0/1504
28	LZ	0.41	0/1130	0.54	0/1507
29	La	0.40	0/1191	0.53	0/1591
30	Lb	0.33	0/889	0.54	0/1175
31	Lc	0.41	0/774	0.58	0/1038
32	Ld	0.39	0/903	0.53	0/1216

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
33	Le	0.42	0/1071	0.57	0/1429
34	Lf	0.46	1/895 (0.1%)	0.61	0/1198
35	Lg	0.40	0/916	0.57	0/1220
36	Lh	0.35	0/1023	0.47	0/1351
37	Li	0.32	0/843	0.46	0/1115
38	Lj	0.43	0/720	0.61	0/952
39	Lk	0.35	0/575	0.53	0/761
40	Ll	0.36	0/454	0.53	0/599
41	Lm	0.37	0/435	0.54	0/575
42	Ln	0.34	0/231	0.43	0/294
43	Lo	0.40	0/876	0.52	0/1156
44	Lp	0.41	0/718	0.54	0/953
45	Lr	0.38	0/1017	0.55	0/1364
46	Ls	0.30	0/1519	0.63	1/2052 (0.0%)
47	Lt	0.35	0/1058	0.82	3/1430 (0.2%)
48	Lz	0.30	0/1769	0.65	1/2371 (0.0%)
49	S2	0.68	1/41245 (0.0%)	1.11	217/64265 (0.3%)
50	SA	0.37	0/1778	0.59	1/2416 (0.0%)
51	SB	0.36	0/1765	0.58	1/2362 (0.0%)
52	SD	0.35	0/1793	0.61	1/2414 (0.0%)
53	SE	0.35	0/2118	0.57	1/2849 (0.0%)
54	SF	0.36	0/1481	0.55	0/1988
55	SH	0.36	0/1519	0.63	1/2033 (0.0%)
56	SI	0.37	0/1715	0.56	0/2287
57	SK	0.31	0/851	0.54	1/1147 (0.1%)
58	SL	0.40	0/1268	0.54	0/1696
59	SP	0.33	0/1082	0.60	1/1446 (0.1%)
60	SQ	0.35	0/1160	0.61	0/1553
61	SR	0.32	0/1105	0.59	1/1484 (0.1%)
62	SS	0.31	0/1216	0.54	0/1628
63	ST	0.33	0/1131	0.56	1/1515 (0.1%)
64	SU	0.32	0/831	0.63	0/1115
65	SV	0.34	0/643	0.56	0/860
66	SX	0.39	0/1116	0.58	0/1490
67	Sa	0.37	0/836	0.56	0/1121
68	Sc	0.33	0/508	0.66	0/680
69	Sd	0.35	0/470	0.55	0/623
70	Sg	0.32	0/2493	0.65	0/3394
71	SC	0.40	0/1762	0.57	0/2381
72	SG	0.32	0/1946	0.59	1/2590 (0.0%)
73	SJ	0.36	0/1550	0.55	0/2069
74	SM	0.32	0/950	0.67	0/1275
75	SN	0.36	0/1232	0.54	1/1656 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
76	SO	0.38	0/1062	0.62	0/1425
77	SW	0.37	0/1051	0.53	0/1406
78	SY	0.34	0/1083	0.53	0/1438
79	SZ	0.32	0/604	0.72	0/810
80	Sb	0.34	0/665	0.55	0/891
81	Se	0.31	0/444	0.55	0/588
82	Sf	0.29	0/560	0.67	0/745
83	CA	0.34	0/2810	0.68	2/3780 (0.1%)
84	CC	0.38	0/1773	1.05	7/2759 (0.3%)
85	CE	0.30	0/1055	0.58	1/1400 (0.1%)
86	CF	0.35	0/244	0.42	0/328
All	All	0.61	3/241703 (0.0%)	0.95	836/354180 (0.2%)

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
1	L5	0	1
4	LA	0	2
5	LB	0	2
11	LH	0	1
13	LJ	0	1
14	LL	0	1
15	LM	0	2
17	LO	0	1
18	LP	0	1
22	LT	0	1
29	La	0	1
34	Lf	0	2
36	Lh	0	1
38	Lj	0	1
46	Ls	0	2
47	Lt	0	3
48	Lz	0	1
51	SB	0	1
52	SD	0	3
54	SF	0	2
55	SH	0	1
56	SI	0	1
60	SQ	0	1

*Continued on next page...*

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Mol	Chain	#Chirality outliers	#Planarity outliers
61	SR	0	1
63	ST	0	1
65	SV	0	1
66	SX	0	3
68	Sc	0	1
73	SJ	0	1
74	SM	0	1
79	SZ	0	2
80	Sb	0	1
All	All	0	45

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
49	S2	1422	G	C6-N1	-6.05	1.35	1.39
34	Lf	47	CYS	CB-SG	-5.20	1.73	1.81
1	L5	4909	A	O3'-P	-5.02	1.55	1.61

All (836) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1417	C	N3-C4-N4	-26.67	99.33	118.00
49	S2	1422	G	N1-C6-O6	-26.29	104.13	119.90
49	S2	1422	G	C5-C6-O6	21.88	141.73	128.60
49	S2	1417	C	C5-C4-N4	20.97	134.88	120.20
49	S2	1772	C	N1-C2-O2	15.42	128.15	118.90
49	S2	1772	C	N3-C2-O2	-15.12	111.32	121.90
1	L5	485	C	C2-N1-C1'	14.57	134.82	118.80
49	S2	501	C	N1-C2-O2	13.24	126.84	118.90
49	S2	501	C	C2-N1-C1'	12.86	132.94	118.80
1	L5	2814	C	N1-C2-O2	12.25	126.25	118.90
49	S2	293	C	N1-C2-O2	12.12	126.17	118.90
1	L5	2710	C	N1-C2-O2	11.82	125.99	118.90
49	S2	882	U	N1-C2-O2	11.60	130.92	122.80
1	L5	3948	C	C2-N1-C1'	11.54	131.49	118.80
1	L5	2710	C	C2-N1-C1'	11.35	131.29	118.80
1	L5	2505	C	N1-C2-O2	11.20	125.62	118.90
1	L5	485	C	C6-N1-C1'	-11.17	107.40	120.80
49	S2	293	C	C2-N1-C1'	11.03	130.94	118.80
1	L5	2814	C	N3-C2-O2	-10.98	114.21	121.90
49	S2	501	C	N3-C2-O2	-10.89	114.28	121.90
49	S2	882	U	C2-N1-C1'	10.86	130.73	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1453	C	N1-C2-O2	10.40	125.14	118.90
49	S2	1453	C	C2-N1-C1'	10.38	130.22	118.80
1	L5	1414	C	N3-C2-O2	-10.34	114.66	121.90
49	S2	1772	C	C6-N1-C2	-10.26	116.20	120.30
1	L5	2710	C	N3-C2-O2	-10.25	114.73	121.90
1	L5	4923	C	N3-C2-O2	-9.98	114.92	121.90
49	S2	118	C	N1-C2-O2	9.80	124.78	118.90
49	S2	1139	C	N3-C2-O2	-9.77	115.06	121.90
1	L5	2505	C	N3-C2-O2	-9.77	115.06	121.90
1	L5	4921	C	N3-C2-O2	-9.74	115.08	121.90
1	L5	485	C	N1-C2-O2	9.69	124.71	118.90
49	S2	882	U	N3-C2-O2	-9.67	115.43	122.20
1	L5	2814	C	C2-N1-C1'	9.65	129.41	118.80
1	L5	100	C	C2-N1-C1'	9.56	129.31	118.80
49	S2	1772	C	C2-N1-C1'	9.54	129.29	118.80
49	S2	1139	C	N1-C2-O2	9.46	124.58	118.90
1	L5	4906	U	N1-C1'-C2'	-9.44	101.61	112.00
1	L5	1414	C	N1-C2-O2	9.38	124.53	118.90
1	L5	4557	U	N3-C2-O2	-9.38	115.64	122.20
1	L5	181	C	N1-C2-O2	9.32	124.49	118.90
1	L5	1367	C	N1-C2-O2	9.32	124.49	118.90
49	S2	1417	C	C4-C5-C6	-9.27	112.77	117.40
1	L5	753	C	N1-C2-O2	9.22	124.43	118.90
1	L5	515	C	N1-C2-O2	9.19	124.42	118.90
49	S2	293	C	N3-C2-O2	-9.11	115.53	121.90
1	L5	4149	C	N3-C2-O2	-9.09	115.53	121.90
49	S2	356	C	C2-N1-C1'	9.07	128.78	118.80
1	L5	753	C	N3-C2-O2	-9.07	115.55	121.90
1	L5	2627	C	C2-N1-C1'	8.97	128.67	118.80
1	L5	417	G	O4'-C1'-N9	8.88	115.30	108.20
49	S2	1417	C	N3-C4-C5	8.87	125.45	121.90
1	L5	4926	C	N1-C2-O2	8.86	124.21	118.90
49	S2	118	C	C2-N1-C1'	8.86	128.54	118.80
49	S2	118	C	N3-C2-O2	-8.81	115.73	121.90
49	S2	501	C	C6-N1-C1'	-8.80	110.24	120.80
49	S2	501	C	C6-N1-C2	-8.80	116.78	120.30
1	L5	1252	C	N3-C2-O2	-8.80	115.74	121.90
1	L5	3948	C	N1-C2-O2	8.70	124.12	118.90
49	S2	1520	G	C4-N9-C1'	8.64	137.73	126.50
1	L5	2410	C	C2-N1-C1'	8.57	128.23	118.80
1	L5	4926	C	C2-N1-C1'	8.55	128.21	118.80
49	S2	1271	C	N1-C2-O2	8.51	124.01	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	100	C	N3-C2-O2	-8.51	115.95	121.90
1	L5	2505	C	C2-N1-C1'	8.50	128.16	118.80
1	L5	4928	C	C2-N1-C1'	8.45	128.09	118.80
1	L5	1367	C	C2-N1-C1'	8.44	128.08	118.80
1	L5	4551	U	N3-C2-O2	-8.42	116.31	122.20
1	L5	2018	C	C5-C6-N1	8.40	125.20	121.00
1	L5	4229	U	N3-C2-O2	-8.37	116.34	122.20
1	L5	233	U	N3-C2-O2	-8.37	116.34	122.20
49	S2	427	U	N3-C2-O2	-8.32	116.37	122.20
1	L5	4557	U	C2-N1-C1'	8.25	127.60	117.70
49	S2	356	C	N1-C2-O2	8.20	123.82	118.90
1	L5	4303	C	N3-C2-O2	-8.18	116.17	121.90
49	S2	1139	C	C2-N1-C1'	8.18	127.79	118.80
1	L5	175	C	N3-C2-O2	-8.13	116.20	121.90
1	L5	1193	C	C2-N1-C1'	8.14	127.75	118.80
3	L8	128	C	C5-C6-N1	8.13	125.07	121.00
1	L5	4101	C	N3-C4-C5	8.12	125.15	121.90
1	L5	925	C	N1-C2-O2	8.12	123.77	118.90
1	L5	4557	U	N1-C2-O2	8.10	128.47	122.80
1	L5	209	U	C2-N1-C1'	8.08	127.39	117.70
1	L5	3948	C	C5-C6-N1	8.06	125.03	121.00
1	L5	5022	U	N1-C2-O2	8.05	128.43	122.80
49	S2	293	C	C6-N1-C1'	-8.04	111.15	120.80
1	L5	1378	C	N1-C2-O2	7.99	123.69	118.90
49	S2	168	C	N1-C2-O2	7.95	123.67	118.90
1	L5	3948	C	C6-N1-C1'	-7.94	111.27	120.80
1	L5	2710	C	C6-N1-C2	-7.93	117.13	120.30
84	CC	74	C	C6-N1-C2	-7.92	117.13	120.30
5	LB	17	LEU	CA-CB-CG	7.92	133.51	115.30
1	L5	2786	C	C6-N1-C2	-7.91	117.14	120.30
3	L8	51	U	N1-C2-O2	7.90	128.33	122.80
3	L8	51	U	N3-C2-O2	-7.88	116.68	122.20
3	L8	64	U	N3-C2-O2	-7.85	116.71	122.20
1	L5	233	U	N1-C2-O2	7.85	128.29	122.80
1	L5	100	C	N1-C2-O2	7.84	123.60	118.90
1	L5	4281	A	N7-C8-N9	7.81	117.70	113.80
1	L5	4758	U	C2-N1-C1'	7.80	127.06	117.70
1	L5	3948	C	C6-N1-C2	-7.79	117.18	120.30
1	L5	2262	G	C4-N9-C1'	7.76	136.58	126.50
49	S2	1520	G	C8-N9-C1'	-7.75	116.92	127.00
1	L5	4303	C	C2-N1-C1'	7.71	127.28	118.80
1	L5	1216	C	C2-N1-C1'	7.71	127.28	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4682	U	N3-C2-O2	-7.70	116.81	122.20
49	S2	1453	C	N3-C2-O2	-7.69	116.52	121.90
1	L5	2710	C	C6-N1-C1'	-7.68	111.58	120.80
49	S2	427	U	C2-N1-C1'	7.68	126.92	117.70
1	L5	3757	G	O4'-C1'-N9	7.67	114.33	108.20
49	S2	179	C	N1-C2-O2	7.66	123.50	118.90
1	L5	4926	C	N3-C2-O2	-7.66	116.54	121.90
1	L5	115	C	N1-C2-O2	7.64	123.48	118.90
1	L5	4758	U	N3-C2-O2	-7.63	116.86	122.20
49	S2	548	C	C2-N1-C1'	7.62	127.19	118.80
1	L5	4758	U	N1-C2-O2	7.62	128.14	122.80
1	L5	181	C	C2-N1-C1'	7.54	127.10	118.80
1	L5	181	C	N3-C2-O2	-7.54	116.62	121.90
1	L5	907	C	C2-N1-C1'	7.54	127.09	118.80
1	L5	1915	C	N3-C2-O2	-7.53	116.63	121.90
1	L5	2505	C	C6-N1-C2	-7.52	117.29	120.30
1	L5	925	C	N3-C2-O2	-7.51	116.64	121.90
1	L5	4910	A	C4'-C3'-O3'	-7.51	93.64	109.40
1	L5	129	C	N3-C2-O2	-7.44	116.69	121.90
1	L5	4303	C	N1-C2-O2	7.44	123.36	118.90
3	L8	128	C	C2-N1-C1'	7.42	126.97	118.80
49	S2	1271	C	N3-C2-O2	-7.41	116.71	121.90
1	L5	4923	C	N1-C2-O2	7.41	123.35	118.90
1	L5	4709	U	N3-C2-O2	-7.41	117.02	122.20
1	L5	4887	C	N1-C2-O2	7.37	123.32	118.90
1	L5	1367	C	N3-C2-O2	-7.37	116.74	121.90
1	L5	2257	C	C2-N1-C1'	7.35	126.88	118.80
1	L5	1714	C	N1-C2-O2	7.31	123.29	118.90
49	S2	427	U	N1-C2-O2	7.31	127.92	122.80
49	S2	1265	A	C2-N3-C4	7.30	114.25	110.60
1	L5	1182	C	N1-C2-O2	7.30	123.28	118.90
49	S2	1022	U	C2-N1-C1'	7.29	126.45	117.70
49	S2	1453	C	C6-N1-C1'	-7.29	112.06	120.80
1	L5	174	C	N3-C2-O2	-7.28	116.81	121.90
1	L5	115	C	C2-N1-C1'	7.28	126.81	118.80
1	L5	963	G	C4-N9-C1'	7.26	135.94	126.50
49	S2	178	C	N1-C2-O2	7.23	123.24	118.90
1	L5	4502	C	N1-C2-O2	7.23	123.24	118.90
1	L5	5022	U	C2-N1-C1'	7.22	126.37	117.70
1	L5	906	C	N3-C2-O2	-7.21	116.86	121.90
1	L5	1703	C	N1-C2-O2	7.20	123.22	118.90
49	S2	630	U	C2-N1-C1'	7.20	126.34	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4709	U	N1-C2-O2	7.20	127.84	122.80
1	L5	4281	A	O4'-C1'-N9	7.19	113.95	108.20
1	L5	516	C	N1-C2-O2	7.19	123.21	118.90
49	S2	882	U	C5-C6-N1	7.19	126.29	122.70
49	S2	1016	U	N3-C2-O2	-7.18	117.17	122.20
1	L5	453	G	N3-C4-C5	-7.17	125.01	128.60
1	L5	3909	C	C6-N1-C2	-7.17	117.43	120.30
1	L5	515	C	C2-N1-C1'	7.17	126.69	118.80
1	L5	2409	U	N1-C2-N3	7.14	119.18	114.90
1	L5	1191	C	N3-C2-O2	-7.14	116.90	121.90
49	S2	882	U	C6-N1-C1'	-7.13	111.22	121.20
1	L5	1340	C	C5-C6-N1	7.13	124.56	121.00
1	L5	3778	U	N1-C2-O2	7.12	127.78	122.80
1	L5	4928	C	N1-C2-O2	7.12	123.17	118.90
1	L5	485	C	C5-C6-N1	7.09	124.54	121.00
1	L5	2262	G	N3-C4-N9	7.08	130.25	126.00
1	L5	1182	C	C2-N1-C1'	7.07	126.57	118.80
1	L5	2528	G	N3-C4-C5	-7.05	125.08	128.60
1	L5	115	C	N3-C2-O2	-7.04	116.97	121.90
1	L5	100	C	C6-N1-C2	-7.04	117.48	120.30
1	L5	2820	C	N1-C2-O2	7.03	123.12	118.90
1	L5	5022	U	N3-C2-O2	-7.02	117.28	122.20
1	L5	2260	C	N1-C2-O2	7.00	123.10	118.90
49	S2	1292	C	N1-C2-O2	6.99	123.10	118.90
1	L5	4281	A	C8-N9-C4	-6.99	103.01	105.80
1	L5	4926	C	C6-N1-C2	-6.97	117.51	120.30
1	L5	1632	A	C2-N3-C4	6.96	114.08	110.60
1	L5	3636	C	C6-N1-C2	-6.96	117.52	120.30
1	L5	4303	C	C6-N1-C2	-6.94	117.53	120.30
1	L5	2018	C	C6-N1-C2	-6.93	117.53	120.30
1	L5	4773	C	N1-C2-O2	6.92	123.05	118.90
49	S2	501	C	C5-C6-N1	6.92	124.46	121.00
1	L5	220	C	C6-N1-C2	-6.91	117.54	120.30
1	L5	499	G	N3-C4-N9	6.90	130.14	126.00
59	SP	107	ILE	C-N-CA	6.90	138.94	121.70
1	L5	41	C	C6-N1-C2	-6.90	117.54	120.30
1	L5	4398	C	N1-C2-O2	6.89	123.03	118.90
1	L5	2528	G	C4-N9-C1'	6.88	135.44	126.50
1	L5	1241	C	N1-C2-O2	6.88	123.03	118.90
1	L5	489	C	C2-N1-C1'	6.87	126.35	118.80
1	L5	220	C	C2-N1-C1'	6.87	126.35	118.80
1	L5	4229	U	N1-C2-O2	6.86	127.60	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	1241	C	C2-N1-C1'	6.85	126.34	118.80
1	L5	4904	G	C4'-C3'-O3'	6.85	126.70	113.00
49	S2	1520	G	N3-C4-N9	6.84	130.11	126.00
1	L5	2627	C	C6-N1-C2	-6.84	117.56	120.30
1	L5	4662	C	C6-N1-C2	-6.83	117.57	120.30
49	S2	1389	C	C2-N1-C1'	6.82	126.30	118.80
1	L5	753	C	C6-N1-C2	-6.81	117.58	120.30
1	L5	2820	C	N3-C2-O2	-6.81	117.14	121.90
1	L5	2260	C	C2-N1-C1'	6.79	126.27	118.80
1	L5	453	G	N3-C4-N9	6.79	130.07	126.00
1	L5	4945	G	N3-C4-N9	6.79	130.07	126.00
49	S2	688	U	P-O3'-C3'	6.78	127.83	119.70
49	S2	1591	C	N1-C2-O2	6.77	122.96	118.90
1	L5	3767	C	C5-C6-N1	6.77	124.39	121.00
1	L5	4924	C	N3-C2-O2	-6.77	117.16	121.90
49	S2	1016	U	N1-C2-O2	6.77	127.54	122.80
1	L5	2410	C	C6-N1-C2	-6.76	117.59	120.30
1	L5	1821	G	N3-C4-N9	6.76	130.06	126.00
1	L5	672	C	C2-N1-C1'	6.75	126.22	118.80
1	L5	3772	U	C2-N1-C1'	6.75	125.80	117.70
1	L5	4928	C	N3-C2-O2	-6.75	117.18	121.90
1	L5	1082	C	OP1-P-O3'	6.74	120.03	105.20
1	L5	499	G	C4-N9-C1'	6.74	135.26	126.50
1	L5	1216	C	N1-C2-O2	6.74	122.94	118.90
1	L5	4682	U	N1-C2-O2	6.74	127.52	122.80
1	L5	914	U	P-O3'-C3'	6.73	127.78	119.70
1	L5	453	G	C4-N9-C1'	6.73	135.24	126.50
1	L5	2760	G	P-O3'-C3'	6.73	127.77	119.70
1	L5	2627	C	N1-C2-O2	6.72	122.93	118.90
1	L5	985	C	C2-N1-C1'	6.70	126.17	118.80
1	L5	1821	G	N3-C4-C5	-6.69	125.26	128.60
1	L5	2262	G	N3-C4-C5	-6.68	125.26	128.60
1	L5	3771	C	C2-N1-C1'	6.67	126.14	118.80
1	L5	209	U	C6-N1-C1'	-6.67	111.86	121.20
1	L5	963	G	C8-N9-C1'	-6.67	118.34	127.00
1	L5	1344	C	C6-N1-C2	-6.66	117.64	120.30
49	S2	579	C	N1-C2-O2	6.66	122.89	118.90
49	S2	494	C	N1-C2-O2	6.65	122.89	118.90
1	L5	2262	G	C8-N9-C1'	-6.65	118.36	127.00
49	S2	356	C	C6-N1-C1'	-6.65	112.82	120.80
1	L5	2351	C	C6-N1-C2	-6.65	117.64	120.30
3	L8	111	U	C2-N1-C1'	6.64	125.67	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	77	U	N3-C2-O2	-6.63	117.56	122.20
1	L5	4502	C	N3-C2-O2	-6.62	117.27	121.90
1	L5	1582	U	N3-C2-O2	-6.61	117.57	122.20
1	L5	1193	C	N1-C2-O2	6.61	122.87	118.90
1	L5	4107	G	N1-C6-O6	-6.61	115.93	119.90
1	L5	516	C	N3-C2-O2	-6.61	117.28	121.90
1	L5	1083	U	O5'-P-OP1	-6.60	99.76	105.70
1	L5	1082	C	P-O3'-C3'	6.60	127.62	119.70
1	L5	1340	C	C6-N1-C2	-6.59	117.66	120.30
1	L5	2814	C	C6-N1-C2	-6.59	117.66	120.30
1	L5	1405	C	N1-C2-O2	6.59	122.85	118.90
49	S2	537	C	N1-C2-O2	6.59	122.85	118.90
49	S2	1453	C	C5-C6-N1	6.58	124.29	121.00
1	L5	3772	U	N3-C2-O2	-6.56	117.61	122.20
1	L5	4887	C	C2-N1-C1'	6.56	126.02	118.80
1	L5	1607	C	C6-N1-C2	-6.56	117.67	120.30
1	L5	2814	C	C6-N1-C1'	-6.56	112.92	120.80
49	S2	1865	C	C6-N1-C2	-6.55	117.68	120.30
1	L5	2675	G	P-O3'-C3'	6.55	127.56	119.70
1	L5	2107	C	N3-C2-O2	-6.53	117.33	121.90
49	S2	659	G	C4-N9-C1'	6.53	134.98	126.50
1	L5	1193	C	C6-N1-C2	-6.52	117.69	120.30
1	L5	2709	C	N3-C4-C5	6.52	124.51	121.90
1	L5	1552	G	O4'-C1'-N9	6.51	113.41	108.20
84	CC	35	U	P-O3'-C3'	6.50	127.50	119.70
49	S2	1453	C	C6-N1-C2	-6.48	117.71	120.30
49	S2	630	U	N1-C2-O2	6.47	127.33	122.80
1	L5	4471	U	N3-C2-O2	-6.47	117.67	122.20
49	S2	130	G	N3-C4-N9	6.47	129.88	126.00
12	LI	212	LEU	CA-CB-CG	6.46	130.17	115.30
49	S2	168	C	N3-C2-O2	-6.46	117.38	121.90
1	L5	3778	U	N3-C2-O2	-6.46	117.68	122.20
1	L5	485	C	C6-N1-C2	-6.46	117.72	120.30
1	L5	1417	C	C2-N1-C1'	6.45	125.89	118.80
49	S2	593	C	N1-C2-O2	6.43	122.76	118.90
49	S2	118	C	C6-N1-C2	-6.42	117.73	120.30
1	L5	1378	C	N3-C2-O2	-6.41	117.41	121.90
49	S2	1597	C	N3-C2-O2	-6.40	117.42	121.90
49	S2	1389	C	C6-N1-C2	-6.40	117.74	120.30
49	S2	1578	U	N1-C2-O2	6.40	127.28	122.80
1	L5	1378	C	C2-N1-C1'	6.39	125.83	118.80
1	L5	1417	C	C5-C6-N1	6.39	124.19	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	2048	U	N1-C2-O2	6.39	127.27	122.80
1	L5	3775	A	N7-C8-N9	6.38	116.99	113.80
1	L5	3762	U	N1-C2-O2	6.38	127.27	122.80
1	L5	1644	C	C6-N1-C2	-6.38	117.75	120.30
1	L5	2802	C	C6-N1-C2	-6.37	117.75	120.30
1	L5	4314	C	N1-C2-O2	6.37	122.72	118.90
1	L5	655	C	N3-C2-O2	-6.37	117.44	121.90
49	S2	179	C	N3-C2-O2	-6.37	117.44	121.90
1	L5	4904	G	C3'-C2'-C1'	6.37	106.59	101.50
49	S2	130	G	C4-N9-C1'	6.36	134.77	126.50
49	S2	1660	C	N3-C2-O2	-6.36	117.45	121.90
1	L5	1663	C	C5-C6-N1	6.34	124.17	121.00
1	L5	4502	C	C6-N1-C2	-6.34	117.77	120.30
1	L5	100	C	C6-N1-C1'	-6.33	113.20	120.80
1	L5	1251	C	N1-C2-O2	6.33	122.70	118.90
1	L5	4945	G	C5-C6-O6	-6.33	124.80	128.60
49	S2	1271	C	C2-N1-C1'	6.33	125.76	118.80
49	S2	1081	U	N3-C2-O2	-6.32	117.77	122.20
1	L5	112	C	C2-N1-C1'	6.32	125.75	118.80
1	L5	4476	C	C2-N1-C1'	6.32	125.75	118.80
1	L5	2409	U	C4-C5-C6	6.32	123.49	119.70
51	SB	62	LEU	CA-CB-CG	6.31	129.81	115.30
1	L5	2257	C	N1-C2-O2	6.31	122.69	118.90
3	L8	111	U	N1-C2-O2	6.30	127.21	122.80
49	S2	325	C	N1-C2-O2	6.30	122.68	118.90
1	L5	3775	A	C2-N3-C4	-6.29	107.45	110.60
1	L5	2464	C	O4'-C1'-N1	6.28	113.23	108.20
1	L5	3892	U	N3-C2-O2	-6.28	117.80	122.20
3	L8	128	C	C6-N1-C2	-6.28	117.79	120.30
49	S2	1420	G	C4-N9-C1'	6.28	134.67	126.50
1	L5	1193	C	C5-C6-N1	6.28	124.14	121.00
52	SD	14	ASP	CB-CG-OD1	6.28	123.95	118.30
1	L5	1417	C	C6-N1-C2	-6.27	117.79	120.30
1	L5	673	C	C2-N1-C1'	6.27	125.70	118.80
1	L5	1714	C	C2-N1-C1'	6.27	125.69	118.80
49	S2	1261	C	N1-C2-O2	6.26	122.66	118.90
49	S2	1315	U	N1-C2-O2	6.26	127.19	122.80
1	L5	3594	C	C6-N1-C2	-6.26	117.80	120.30
1	L5	963	G	N3-C4-N9	6.26	129.75	126.00
49	S2	1139	C	C6-N1-C2	-6.25	117.80	120.30
1	L5	1216	C	N3-C2-O2	-6.25	117.53	121.90
53	SE	139	LEU	CA-CB-CG	6.25	129.66	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
85	CE	199	LEU	CA-CB-CG	6.24	129.66	115.30
1	L5	499	G	N3-C4-C5	-6.23	125.49	128.60
1	L5	456	C	O4'-C1'-N1	6.23	113.18	108.20
1	L5	2033	A	P-O3'-C3'	6.23	127.17	119.70
3	L8	64	U	N1-C2-O2	6.21	127.15	122.80
2	L7	28	C	C6-N1-C2	-6.21	117.82	120.30
49	S2	356	C	N3-C2-O2	-6.18	117.57	121.90
1	L5	4360	U	N3-C2-O2	-6.18	117.87	122.20
49	S2	1578	U	N3-C2-O2	-6.17	117.88	122.20
1	L5	4149	C	N1-C2-O2	6.17	122.60	118.90
1	L5	4928	C	C6-N1-C1'	-6.16	113.41	120.80
1	L5	3767	C	C6-N1-C2	-6.14	117.84	120.30
1	L5	1182	C	N3-C2-O2	-6.14	117.60	121.90
1	L5	2094	G	C4-N9-C1'	6.13	134.47	126.50
1	L5	2416	G	P-O3'-C3'	6.12	127.05	119.70
49	S2	1660	C	C2-N1-C1'	6.12	125.53	118.80
49	S2	325	C	C2-N1-C1'	6.11	125.52	118.80
1	L5	4913	G	P-O3'-C3'	6.10	127.02	119.70
1	L5	41	C	C5-C6-N1	6.10	124.05	121.00
1	L5	4747	C	C2-N1-C1'	6.10	125.50	118.80
49	S2	1309	C	C2-N1-C1'	6.08	125.48	118.80
49	S2	1865	C	N3-C2-O2	-6.08	117.65	121.90
1	L5	1582	U	N1-C2-O2	6.08	127.05	122.80
1	L5	2107	C	N1-C2-O2	6.07	122.54	118.90
49	S2	1660	C	N1-C2-O2	6.07	122.54	118.90
49	S2	632	C	C6-N1-C2	-6.07	117.87	120.30
49	S2	548	C	N1-C2-O2	6.07	122.54	118.90
1	L5	2627	C	C5-C6-N1	6.06	124.03	121.00
1	L5	3909	C	N3-C2-O2	-6.06	117.66	121.90
49	S2	824	C	N1-C2-O2	6.05	122.53	118.90
49	S2	1292	C	N3-C2-O2	-6.05	117.67	121.90
47	Lt	121	LEU	CA-CB-CG	6.05	129.21	115.30
1	L5	1977	C	P-O3'-C3'	6.04	126.95	119.70
49	S2	1420	G	N3-C4-C5	-6.04	125.58	128.60
49	S2	130	G	N3-C4-C5	-6.03	125.58	128.60
49	S2	179	C	C2-N1-C1'	6.03	125.43	118.80
49	S2	1865	C	C2-N1-C1'	6.03	125.43	118.80
1	L5	1458	C	N1-C2-O2	6.01	122.51	118.90
1	L5	3641	U	C4-C5-C6	6.01	123.31	119.70
1	L5	2011	C	N1-C2-O2	6.01	122.50	118.90
2	L7	39	C	N1-C2-O2	6.01	122.50	118.90
49	S2	4	C	C6-N1-C2	-6.00	117.90	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4107	G	C4-N9-C1'	5.99	134.29	126.50
49	S2	1420	G	N3-C4-N9	5.99	129.59	126.00
1	L5	1666	C	C6-N1-C2	-5.99	117.90	120.30
49	S2	178	C	N3-C2-O2	-5.99	117.71	121.90
1	L5	499	G	C8-N9-C1'	-5.99	119.22	127.00
49	S2	293	C	C6-N1-C2	-5.99	117.91	120.30
1	L5	925	C	C6-N1-C2	-5.98	117.91	120.30
1	L5	2096	G	C4-N9-C1'	5.98	134.27	126.50
49	S2	310	C	C2-N1-C1'	5.97	125.37	118.80
49	S2	1022	U	N3-C2-O2	-5.97	118.02	122.20
1	L5	4880	C	N1-C2-O2	5.97	122.48	118.90
1	L5	4398	C	N3-C2-O2	-5.97	117.72	121.90
1	L5	1241	C	N3-C2-O2	-5.97	117.72	121.90
49	S2	814	U	N3-C2-O2	-5.97	118.02	122.20
1	L5	2410	C	C5-C6-N1	5.96	123.98	121.00
1	L5	2627	C	C6-N1-C1'	-5.96	113.64	120.80
2	L7	102	U	N1-C2-O2	5.96	126.97	122.80
49	S2	930	C	N1-C2-O2	5.95	122.47	118.90
1	L5	906	C	C6-N1-C2	-5.95	117.92	120.30
1	L5	2112	G	C1'-C2'-O2'	-5.95	92.76	110.60
1	L5	2708	U	C2-N1-C1'	5.94	124.83	117.70
1	L5	4773	C	N3-C2-O2	-5.94	117.74	121.90
1	L5	2351	C	C2-N1-C1'	5.94	125.33	118.80
1	L5	1853	G	C4-N9-C1'	5.94	134.22	126.50
49	S2	632	C	C2-N1-C1'	5.94	125.33	118.80
49	S2	119	U	N3-C2-O2	-5.93	118.05	122.20
1	L5	1447	C	C2-N1-C1'	5.93	125.33	118.80
1	L5	3685	C	C6-N1-C2	-5.93	117.93	120.30
49	S2	537	C	C2-N1-C1'	5.93	125.33	118.80
49	S2	563	G	P-O3'-C3'	5.93	126.82	119.70
49	S2	1597	C	N1-C2-O2	5.93	122.46	118.90
1	L5	2409	U	N3-C2-O2	-5.93	118.05	122.20
49	S2	1434	C	P-O3'-C3'	5.92	126.81	119.70
1	L5	233	U	C2-N1-C1'	5.92	124.81	117.70
49	S2	1772	C	C5-C6-N1	5.92	123.96	121.00
1	L5	485	C	N3-C2-O2	-5.92	117.76	121.90
49	S2	119	U	N1-C2-O2	5.92	126.94	122.80
1	L5	2528	G	N3-C4-N9	5.92	129.55	126.00
1	L5	4241	C	C2-N1-C1'	5.91	125.31	118.80
1	L5	4352	U	N3-C2-O2	-5.91	118.06	122.20
49	S2	118	C	C6-N1-C1'	-5.91	113.70	120.80
1	L5	1458	C	N3-C2-O2	-5.91	117.76	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4887	C	N3-C2-O2	-5.91	117.76	121.90
1	L5	2351	C	C5-C6-N1	5.90	123.95	121.00
1	L5	4314	C	N3-C2-O2	-5.90	117.77	121.90
49	S2	168	C	C6-N1-C2	-5.90	117.94	120.30
49	S2	674	C	C6-N1-C2	-5.90	117.94	120.30
49	S2	1590	C	C6-N1-C2	-5.90	117.94	120.30
49	S2	1047	C	C6-N1-C2	-5.89	117.94	120.30
49	S2	1218	C	C5-C6-N1	5.89	123.94	121.00
3	L8	111	U	N3-C2-O2	-5.89	118.08	122.20
49	S2	314	U	N3-C2-O2	-5.89	118.08	122.20
1	L5	2426	U	N3-C2-O2	-5.88	118.08	122.20
1	L5	4921	C	N1-C2-O2	5.87	122.42	118.90
2	L7	29	C	C2-N1-C1'	5.87	125.26	118.80
49	S2	688	U	N3-C2-O2	-5.87	118.09	122.20
1	L5	282	C	N1-C2-O2	5.86	122.42	118.90
1	L5	205	C	N3-C2-O2	-5.86	117.80	121.90
1	L5	1082	C	O4'-C1'-N1	5.86	112.88	108.20
1	L5	1367	C	C6-N1-C1'	-5.85	113.78	120.80
49	S2	1821	U	N3-C2-O2	-5.85	118.11	122.20
1	L5	1755	C	N1-C2-O2	5.85	122.41	118.90
1	L5	4171	C	N1-C2-O2	5.84	122.41	118.90
49	S2	112	U	P-O3'-C3'	5.84	126.71	119.70
1	L5	2394	G	O4'-C1'-N9	5.84	112.87	108.20
1	L5	1607	C	N3-C2-O2	-5.84	117.81	121.90
1	L5	1703	C	N3-C2-O2	-5.84	117.81	121.90
49	S2	1117	C	N1-C2-O2	5.83	122.40	118.90
49	S2	659	G	C8-N9-C1'	-5.83	119.42	127.00
49	S2	1216	C	N1-C2-O2	5.82	122.39	118.90
1	L5	5028	G	N3-C4-N9	5.82	129.49	126.00
1	L5	3948	C	N3-C2-O2	-5.82	117.83	121.90
1	L5	220	C	C5-C6-N1	5.82	123.91	121.00
49	S2	1315	U	N3-C2-O2	-5.82	118.13	122.20
1	L5	2018	C	C2-N1-C1'	5.82	125.20	118.80
49	S2	814	U	N1-C2-O2	5.82	126.87	122.80
1	L5	2860	C	C6-N1-C2	-5.82	117.97	120.30
1	L5	453	G	C2-N3-C4	5.81	114.81	111.90
5	LB	360	LEU	CA-CB-CG	5.81	128.67	115.30
1	L5	2255	C	C2-N1-C1'	5.81	125.19	118.80
1	L5	5035	U	N3-C2-O2	-5.81	118.14	122.20
84	CC	74	C	P-O3'-C3'	5.81	126.67	119.70
83	CA	156	LEU	CA-CB-CG	5.80	128.64	115.30
2	L7	102	U	N3-C2-O2	-5.80	118.14	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1271	C	C6-N1-C2	-5.80	117.98	120.30
1	L5	2048	U	N3-C2-O2	-5.79	118.14	122.20
49	S2	402	C	C6-N1-C2	-5.79	117.98	120.30
1	L5	515	C	C5-C6-N1	5.79	123.89	121.00
1	L5	1472	C	C2-N1-C1'	5.79	125.16	118.80
1	L5	1663	C	C2-N1-C1'	5.79	125.16	118.80
1	L5	3967	G	N3-C4-C5	-5.79	125.71	128.60
49	S2	1219	C	N1-C2-O2	5.79	122.37	118.90
49	S2	1022	U	N1-C2-O2	5.78	126.85	122.80
1	L5	3762	U	N3-C2-O2	-5.78	118.15	122.20
1	L5	181	C	C6-N1-C2	-5.78	117.99	120.30
1	L5	4392	G	N1-C6-O6	-5.78	116.44	119.90
1	L5	1447	C	C6-N1-C2	-5.77	117.99	120.30
1	L5	2667	C	N1-C2-O2	5.76	122.36	118.90
1	L5	515	C	N3-C2-O2	-5.76	117.87	121.90
1	L5	4594	U	N3-C2-O2	-5.76	118.17	122.20
2	L7	28	C	N1-C2-O2	5.76	122.36	118.90
1	L5	1893	C	C6-N1-C2	-5.76	118.00	120.30
1	L5	1915	C	N1-C2-O2	5.76	122.35	118.90
1	L5	3772	U	N1-C2-O2	5.75	126.83	122.80
5	LB	17	LEU	CB-CG-CD2	-5.75	101.23	111.00
1	L5	4864	U	N1-C2-O2	5.74	126.82	122.80
1	L5	1405	C	N3-C2-O2	-5.73	117.89	121.90
1	L5	1602	U	N3-C2-O2	-5.73	118.19	122.20
1	L5	4293	U	C2-N1-C1'	5.73	124.58	117.70
49	S2	331	C	C2-N1-C1'	5.73	125.10	118.80
1	L5	100	C	O4'-C1'-N1	5.73	112.78	108.20
1	L5	2362	U	N3-C2-O2	-5.73	118.19	122.20
1	L5	1762	C	C6-N1-C2	-5.73	118.01	120.30
1	L5	1439	C	C2-N1-C1'	5.72	125.10	118.80
1	L5	3636	C	N3-C2-O2	-5.72	117.89	121.90
1	L5	3882	C	C2-N1-C1'	5.72	125.09	118.80
1	L5	294	G	C4-N9-C1'	5.71	133.93	126.50
49	S2	291	G	P-O3'-C3'	5.71	126.55	119.70
49	S2	579	C	N3-C2-O2	-5.71	117.91	121.90
49	S2	1116	C	N1-C2-O2	5.71	122.32	118.90
1	L5	2786	C	P-O3'-C3'	5.70	126.54	119.70
1	L5	655	C	N1-C2-O2	5.70	122.32	118.90
1	L5	4101	C	N3-C4-N4	-5.70	114.01	118.00
1	L5	4594	U	C2-N1-C1'	5.70	124.54	117.70
1	L5	2867	C	C6-N1-C2	-5.70	118.02	120.30
1	L5	2529	A	O4'-C1'-N9	5.70	112.76	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	2860	C	N3-C2-O2	-5.70	117.91	121.90
49	S2	1261	C	C6-N1-C2	-5.69	118.02	120.30
1	L5	4904	G	N9-C1'-C2'	-5.69	105.74	112.00
49	S2	1520	G	N3-C4-C5	-5.69	125.76	128.60
3	L8	51	U	C2-N1-C1'	5.68	124.52	117.70
1	L5	1633	G	P-O3'-C3'	5.68	126.52	119.70
1	L5	4281	A	N1-C2-N3	5.67	132.14	129.30
14	LL	144	LEU	CA-CB-CG	5.67	128.35	115.30
1	L5	1809	C	C2-N1-C1'	5.67	125.04	118.80
49	S2	578	C	N1-C2-O2	5.67	122.30	118.90
1	L5	5016	A	O4'-C1'-N9	5.67	112.73	108.20
46	Ls	47	LEU	CA-CB-CG	5.67	128.34	115.30
1	L5	4921	C	C6-N1-C2	-5.67	118.03	120.30
72	SG	41	LEU	CA-CB-CG	5.66	128.32	115.30
1	L5	2806	A	O4'-C1'-N9	5.66	112.72	108.20
1	L5	3778	U	C2-N1-C1'	5.65	124.48	117.70
49	S2	293	C	C5-C6-N1	5.65	123.83	121.00
1	L5	4258	C	C5-C6-N1	5.65	123.83	121.00
1	L5	1309	C	C6-N1-C2	-5.65	118.04	120.30
1	L5	449	C	N1-C2-O2	5.65	122.29	118.90
1	L5	904	C	C6-N1-C2	-5.64	118.04	120.30
1	L5	1252	C	C6-N1-C2	-5.64	118.04	120.30
1	L5	1308	C	C6-N1-C2	-5.64	118.05	120.30
1	L5	2710	C	C5-C6-N1	5.63	123.82	121.00
1	L5	3673	C	P-O3'-C3'	5.63	126.46	119.70
1	L5	2410	C	C6-N1-C1'	-5.63	114.05	120.80
1	L5	4254	G	N3-C4-C5	-5.63	125.79	128.60
1	L5	4281	A	C5-N7-C8	-5.62	101.09	103.90
49	S2	494	C	N3-C2-O2	-5.62	117.96	121.90
49	S2	632	C	C5-C6-N1	5.62	123.81	121.00
1	L5	4923	C	C6-N1-C2	-5.62	118.05	120.30
1	L5	904	C	N1-C2-O2	5.61	122.27	118.90
1	L5	1520	C	O5'-P-OP2	-5.61	100.65	105.70
50	SA	206	ASP	CB-CG-OD1	5.61	123.35	118.30
1	L5	2018	C	N1-C2-O2	5.59	122.25	118.90
1	L5	242	U	N3-C2-O2	-5.59	118.29	122.20
1	L5	1714	C	N3-C2-O2	-5.59	117.99	121.90
1	L5	3618	C	C6-N1-C2	-5.58	118.07	120.30
49	S2	1417	C	N1-C2-N3	-5.58	115.30	119.20
1	L5	195	C	C6-N1-C2	-5.57	118.07	120.30
49	S2	630	U	N3-C2-O2	-5.57	118.30	122.20
1	L5	3657	U	N3-C2-O2	-5.57	118.30	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1219	C	C2-N1-C1'	5.57	124.92	118.80
49	S2	1261	C	C2-N1-C1'	5.56	124.91	118.80
1	L5	2260	C	N3-C2-O2	-5.55	118.01	121.90
1	L5	4969	C	C6-N1-C2	-5.55	118.08	120.30
49	S2	1415	C	C2-N1-C1'	5.55	124.91	118.80
49	S2	542	U	N1-C2-O2	5.55	126.69	122.80
1	L5	282	C	N3-C2-O2	-5.55	118.01	121.90
1	L5	1245	C	C2-N1-C1'	5.55	124.91	118.80
1	L5	3641	U	C5-C6-N1	-5.55	119.92	122.70
1	L5	4600	G	O4'-C1'-N9	5.55	112.64	108.20
49	S2	402	C	C5-C6-N1	5.55	123.77	121.00
49	S2	1016	U	C2-N1-C1'	5.55	124.36	117.70
1	L5	4243	C	C6-N1-C2	-5.55	118.08	120.30
63	ST	110	LEU	CA-CB-CG	5.55	128.06	115.30
1	L5	2416	G	OP2-P-O3'	5.54	117.40	105.20
49	S2	1821	U	N1-C2-O2	5.54	126.68	122.80
1	L5	294	G	N3-C4-C5	-5.54	125.83	128.60
1	L5	4471	U	N1-C2-O2	5.54	126.68	122.80
1	L5	4973	U	N3-C2-O2	-5.54	118.32	122.20
1	L5	453	G	C8-N9-C1'	-5.54	119.80	127.00
1	L5	985	C	C6-N1-C2	-5.54	118.08	120.30
1	L5	4920	C	N1-C2-O2	5.53	122.22	118.90
1	L5	4885	U	N3-C2-O2	-5.53	118.33	122.20
1	L5	907	C	N1-C2-O2	5.53	122.22	118.90
1	L5	985	C	N1-C2-O2	5.52	122.22	118.90
1	L5	228	C	C6-N1-C2	-5.52	118.09	120.30
1	L5	2410	C	N1-C2-O2	5.52	122.21	118.90
1	L5	2900	U	C2-N1-C1'	5.52	124.32	117.70
49	S2	4	C	C2-N1-C1'	5.52	124.87	118.80
1	L5	4293	U	N3-C2-O2	-5.52	118.34	122.20
1	L5	4709	U	C2-N1-C1'	5.51	124.32	117.70
49	S2	1013	U	N3-C2-O2	-5.51	118.34	122.20
49	S2	592	C	C2-N1-C1'	5.51	124.86	118.80
49	S2	1315	U	C2-N1-C1'	5.51	124.31	117.70
1	L5	1663	C	C6-N1-C2	-5.51	118.10	120.30
1	L5	4775	C	C2-N1-C1'	5.51	124.86	118.80
1	L5	4775	C	N1-C2-O2	5.50	122.20	118.90
49	S2	1314	U	C2-N1-C1'	5.50	124.30	117.70
1	L5	1821	G	C4-N9-C1'	5.50	133.65	126.50
1	L5	2899	C	C2-N1-C1'	5.50	124.85	118.80
49	S2	130	G	C8-N9-C1'	-5.50	119.86	127.00
1	L5	1893	C	C2-N1-C1'	5.49	124.84	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	2783	A	N1-C2-N3	-5.49	126.55	129.30
84	CC	74	C	C2'-C3'-O3'	5.49	122.49	113.70
1	L5	4476	C	N1-C2-O2	5.49	122.19	118.90
49	S2	323	C	N1-C2-O2	5.49	122.19	118.90
1	L5	209	U	N1-C2-O2	5.49	126.64	122.80
1	L5	4254	G	C4-N9-C1'	5.48	133.63	126.50
1	L5	4162	C	N1-C2-O2	5.48	122.19	118.90
1	L5	516	C	C2-N1-C1'	5.48	124.83	118.80
1	L5	1216	C	C6-N1-C2	-5.48	118.11	120.30
2	L7	34	C	N1-C2-O2	5.47	122.19	118.90
1	L5	1987	C	C6-N1-C2	-5.47	118.11	120.30
1	L5	2856	C	N1-C2-O2	5.47	122.18	118.90
1	L5	4926	C	C6-N1-C1'	-5.47	114.23	120.80
1	L5	4594	U	N1-C2-O2	5.47	126.63	122.80
1	L5	4945	G	C4-C5-N7	5.47	112.99	110.80
8	LE	50	LEU	CA-CB-CG	5.47	127.88	115.30
1	L5	925	C	C5-C6-N1	5.46	123.73	121.00
1	L5	4107	G	C5-C6-O6	5.46	131.88	128.60
49	S2	168	C	C2-N1-C1'	5.46	124.81	118.80
1	L5	2850	A	C2-N3-C4	5.46	113.33	110.60
1	L5	4450	U	N3-C2-O2	-5.46	118.38	122.20
1	L5	4149	C	C6-N1-C2	-5.46	118.12	120.30
49	S2	578	C	N3-C2-O2	-5.46	118.08	121.90
1	L5	985	C	C5-C6-N1	5.45	123.73	121.00
1	L5	4107	G	N3-C4-C5	-5.45	125.87	128.60
1	L5	4895	C	N1-C2-O2	5.45	122.17	118.90
49	S2	1261	C	N3-C2-O2	-5.45	118.08	121.90
1	L5	2257	C	C6-N1-C1'	-5.44	114.27	120.80
1	L5	4162	C	N3-C2-O2	-5.44	118.09	121.90
49	S2	179	C	C6-N1-C2	-5.44	118.12	120.30
1	L5	4215	C	N1-C2-O2	5.44	122.16	118.90
57	SK	83	LEU	CA-CB-CG	5.44	127.81	115.30
1	L5	1929	A	C4-N9-C1'	5.43	136.08	126.30
49	S2	824	C	N3-C2-O2	-5.43	118.10	121.90
49	S2	1591	C	N3-C2-O2	-5.43	118.10	121.90
49	S2	1078	C	C2-N1-C1'	5.43	124.77	118.80
1	L5	3771	C	N1-C2-O2	5.42	122.15	118.90
1	L5	1251	C	N3-C2-O2	-5.42	118.11	121.90
1	L5	3648	A	O4'-C1'-N9	5.42	112.54	108.20
1	L5	3771	C	C6-N1-C2	-5.42	118.13	120.30
1	L5	4557	U	C6-N1-C1'	-5.42	113.61	121.20
1	L5	3866	C	C6-N1-C2	-5.41	118.14	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	5028	G	N3-C4-C5	-5.41	125.90	128.60
49	S2	1649	U	N1-C2-O2	5.41	126.59	122.80
49	S2	877	C	C2-N1-C1'	5.40	124.75	118.80
1	L5	4398	C	C6-N1-C2	-5.40	118.14	120.30
1	L5	1367	C	C6-N1-C2	-5.39	118.14	120.30
49	S2	950	C	C2-N1-C1'	5.39	124.73	118.80
1	L5	963	G	N3-C4-C5	-5.38	125.91	128.60
1	L5	4450	U	C2-N1-C1'	5.38	124.16	117.70
1	L5	3775	A	C5-N7-C8	-5.38	101.21	103.90
1	L5	4683	U	N3-C2-O2	-5.38	118.44	122.20
1	L5	3680	U	N1-C2-O2	5.37	126.56	122.80
61	SR	109	LEU	CA-CB-CG	5.37	127.64	115.30
49	S2	1265	A	N3-C4-N9	5.36	131.69	127.40
49	S2	950	C	C6-N1-C2	-5.36	118.16	120.30
1	L5	1344	C	C5-C6-N1	5.36	123.68	121.00
1	L5	3594	C	C5-C6-N1	5.36	123.68	121.00
1	L5	4314	C	C6-N1-C2	-5.36	118.16	120.30
1	L5	3841	C	C6-N1-C2	-5.35	118.16	120.30
49	S2	1139	C	C6-N1-C1'	-5.35	114.38	120.80
1	L5	1633	G	OP1-P-O3'	5.34	116.96	105.20
1	L5	4600	G	C4-N9-C1'	-5.34	119.55	126.50
49	S2	1556	A	C2-N3-C4	5.34	113.27	110.60
75	SN	80	LEU	CA-CB-CG	5.34	127.59	115.30
1	L5	5016	A	C5-N7-C8	-5.34	101.23	103.90
1	L5	3939	G	N3-C4-C5	-5.34	125.93	128.60
1	L5	739	G	N3-C4-N9	5.34	129.20	126.00
1	L5	4420	U	N1-C2-O2	5.34	126.54	122.80
2	L7	28	C	N3-C2-O2	-5.34	118.16	121.90
1	L5	2528	G	C8-N9-C1'	-5.33	120.06	127.00
1	L5	984	C	C2-N1-C1'	5.33	124.67	118.80
49	S2	548	C	C6-N1-C1'	-5.33	114.40	120.80
1	L5	1929	A	C2-N3-C4	5.33	113.27	110.60
1	L5	516	C	C6-N1-C2	-5.33	118.17	120.30
1	L5	2255	C	N1-C2-O2	5.33	122.10	118.90
1	L5	4612	C	N1-C2-O2	5.33	122.10	118.90
1	L5	2532	C	C5-C6-N1	5.33	123.66	121.00
1	L5	673	C	C5-C6-N1	5.32	123.66	121.00
49	S2	1865	C	N1-C2-O2	5.32	122.09	118.90
1	L5	4420	U	N3-C2-O2	-5.32	118.48	122.20
1	L5	1579	C	C6-N1-C2	-5.32	118.17	120.30
1	L5	2528	G	C2-N3-C4	5.31	114.56	111.90
3	L8	99	U	N3-C2-O2	-5.31	118.48	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	1252	C	N1-C2-O2	5.31	122.08	118.90
49	S2	1556	A	N3-C4-N9	5.31	131.65	127.40
49	S2	1415	C	C6-N1-C2	-5.30	118.18	120.30
1	L5	1182	C	C6-N1-C2	-5.30	118.18	120.30
49	S2	1591	C	C2-N1-C1'	5.30	124.63	118.80
49	S2	1218	C	C6-N1-C2	-5.30	118.18	120.30
1	L5	205	C	N1-C2-O2	5.30	122.08	118.90
1	L5	972	C	N1-C2-O2	5.30	122.08	118.90
1	L5	4864	U	N3-C2-O2	-5.30	118.49	122.20
1	L5	1726	U	N3-C2-O2	-5.29	118.49	122.20
2	L7	29	C	C6-N1-C2	-5.29	118.18	120.30
1	L5	282	C	C6-N1-C2	-5.29	118.18	120.30
1	L5	4973	U	N1-C2-O2	5.29	126.50	122.80
1	L5	36	U	N3-C2-O2	-5.29	118.50	122.20
49	S2	1259	A	C2-N3-C4	5.29	113.24	110.60
49	S2	1772	C	C6-N1-C1'	-5.28	114.46	120.80
1	L5	4703	U	N3-C2-O2	-5.28	118.50	122.20
1	L5	3775	A	C8-N9-C4	-5.28	103.69	105.80
1	L5	4714	C	N1-C2-O2	5.28	122.07	118.90
1	L5	4926	C	C5-C6-N1	5.28	123.64	121.00
1	L5	1816	C	C6-N1-C2	-5.28	118.19	120.30
1	L5	2409	U	C5-C6-N1	-5.28	120.06	122.70
1	L5	1720	C	C6-N1-C2	-5.27	118.19	120.30
49	S2	1556	A	C4-N9-C1'	5.27	135.79	126.30
1	L5	1395	U	N3-C2-O2	-5.27	118.51	122.20
1	L5	2505	C	C6-N1-C1'	-5.27	114.48	120.80
1	L5	3631	U	N3-C2-O2	-5.27	118.51	122.20
1	L5	4360	U	N1-C2-O2	5.26	126.48	122.80
1	L5	1644	C	C5-C6-N1	5.26	123.63	121.00
1	L5	4476	C	N3-C2-O2	-5.26	118.22	121.90
1	L5	4887	C	C6-N1-C2	-5.26	118.20	120.30
1	L5	96	U	N3-C2-O2	-5.26	118.52	122.20
1	L5	3693	U	N3-C2-O2	-5.26	118.52	122.20
1	L5	757	G	N3-C4-N9	5.25	129.15	126.00
1	L5	673	C	C6-N1-C2	-5.25	118.20	120.30
1	L5	4885	U	N1-C2-O2	5.25	126.47	122.80
1	L5	3637	U	N3-C2-O2	-5.25	118.53	122.20
1	L5	2528	G	C8-N9-C4	-5.24	104.30	106.40
49	S2	1420	G	C8-N9-C1'	-5.24	120.19	127.00
84	CC	74	C	C6-N1-C1'	5.24	127.08	120.80
1	L5	1702	C	C2-N1-C1'	5.24	124.56	118.80
1	L5	2900	U	N1-C2-O2	5.24	126.47	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	2532	C	C6-N1-C2	-5.24	118.21	120.30
1	L5	3930	U	N1-C2-O2	5.23	126.46	122.80
1	L5	5016	A	N7-C8-N9	5.23	116.42	113.80
49	S2	930	C	N3-C2-O2	-5.23	118.24	121.90
55	SH	35	ASP	CB-CG-OD1	5.23	123.01	118.30
1	L5	753	C	C2-N1-C1'	5.22	124.55	118.80
1	L5	1193	C	C6-N1-C1'	-5.22	114.53	120.80
1	L5	5028	G	C4-N9-C1'	5.22	133.29	126.50
49	S2	630	U	C6-N1-C1'	-5.22	113.89	121.20
83	CA	48	LEU	CA-CB-CG	5.22	127.30	115.30
2	L7	109	U	OP2-P-O3'	5.22	116.68	105.20
3	L8	101	C	C6-N1-C2	-5.22	118.21	120.30
7	LD	235	MET	CA-CB-CG	5.21	122.17	113.30
1	L5	294	G	N3-C4-N9	5.21	129.13	126.00
49	S2	1520	G	C6-C5-N7	-5.21	127.27	130.40
1	L5	904	C	C5-C6-N1	5.21	123.61	121.00
1	L5	2018	C	C2-N3-C4	5.20	122.50	119.90
1	L5	1216	C	C6-N1-C1'	-5.19	114.57	120.80
1	L5	4254	G	N3-C4-N9	5.19	129.12	126.00
1	L5	1276	C	C6-N1-C2	-5.19	118.22	120.30
1	L5	3584	C	N1-C2-O2	5.19	122.01	118.90
1	L5	406	C	P-O3'-C3'	5.19	125.92	119.70
1	L5	2304	U	N3-C2-O2	-5.19	118.57	122.20
1	L5	2856	C	N3-C2-O2	-5.19	118.27	121.90
1	L5	3762	U	C2-N1-C1'	5.19	123.92	117.70
1	L5	3967	G	C4-N9-C1'	5.18	133.24	126.50
1	L5	4764	A	N1-C2-N3	-5.18	126.71	129.30
1	L5	907	C	C6-N1-C1'	-5.18	114.58	120.80
1	L5	4206	C	C2-N1-C1'	5.18	124.50	118.80
1	L5	1726	U	N1-C2-O2	5.18	126.42	122.80
49	S2	1389	C	C5-C6-N1	5.18	123.59	121.00
1	L5	3775	A	N1-C2-N3	5.17	131.89	129.30
1	L5	4399	U	N3-C2-O2	-5.17	118.58	122.20
3	L8	128	C	N1-C2-O2	5.17	122.00	118.90
1	L5	2632	U	N3-C2-O2	-5.17	118.58	122.20
1	L5	2532	C	C2-N1-C1'	5.16	124.48	118.80
49	S2	1578	U	C2-N1-C1'	5.16	123.89	117.70
49	S2	1306	U	C2-N1-C1'	5.16	123.89	117.70
1	L5	1735	U	N1-C2-O2	5.16	126.41	122.80
1	L5	1893	C	C5-C6-N1	5.15	123.58	121.00
1	L5	4945	G	N9-C4-C5	-5.15	103.34	105.40
49	S2	418	A	O5'-P-OP1	-5.15	101.06	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	S2	1408	U	C2-N1-C1'	5.15	123.88	117.70
49	S2	325	C	N3-C2-O2	-5.15	118.30	121.90
1	L5	1517	G	N3-C4-N9	5.14	129.09	126.00
1	L5	4901	G	N1-C6-O6	-5.14	116.81	119.90
49	S2	1649	U	N3-C2-O2	-5.14	118.60	122.20
1	L5	907	C	C6-N1-C2	-5.14	118.24	120.30
1	L5	1597	G	O4'-C1'-N9	5.14	112.31	108.20
1	L5	4708	A	C4-N9-C1'	5.14	135.55	126.30
1	L5	515	C	C6-N1-C2	-5.13	118.25	120.30
47	Lt	62	LEU	CA-CB-CG	5.13	127.11	115.30
49	S2	17	C	C6-N1-C2	-5.13	118.25	120.30
1	L5	2371	U	N3-C2-O2	-5.13	118.61	122.20
49	S2	1358	U	N3-C2-O2	-5.13	118.61	122.20
1	L5	3638	G	C2-N3-C4	5.12	114.46	111.90
1	L5	2632	U	N1-C2-O2	5.12	126.38	122.80
49	S2	1149	A	C2-N3-C4	5.12	113.16	110.60
1	L5	4714	C	N3-C2-O2	-5.12	118.32	121.90
47	Lt	37	LEU	CA-CB-CG	5.12	127.06	115.30
1	L5	1439	C	C6-N1-C2	-5.11	118.25	120.30
1	L5	1173	G	N3-C4-N9	-5.11	122.93	126.00
2	L7	39	C	N3-C2-O2	-5.11	118.32	121.90
15	LM	87	ALA	C-N-CA	5.11	134.47	121.70
49	S2	1060	A	O4'-C1'-N9	5.11	112.29	108.20
49	S2	1422	G	C5-C6-N1	5.11	114.06	111.50
1	L5	4093	G	C4-N9-C1'	5.11	133.14	126.50
1	L5	1259	G	N1-C6-O6	-5.11	116.84	119.90
1	L5	4898	G	N1-C6-O6	-5.11	116.84	119.90
1	L5	2802	C	C5-C6-N1	5.10	123.55	121.00
49	S2	1696	C	N1-C2-O2	5.10	121.96	118.90
1	L5	2096	G	C8-N9-C1'	-5.10	120.37	127.00
1	L5	322	C	C6-N1-C2	-5.10	118.26	120.30
1	L5	180	C	C5-C6-N1	5.10	123.55	121.00
1	L5	2867	C	C2-N1-C1'	5.10	124.41	118.80
1	L5	2627	C	N3-C2-O2	-5.10	118.33	121.90
1	L5	4880	C	N3-C2-O2	-5.10	118.33	121.90
49	S2	585	C	C2-N1-C1'	5.10	124.41	118.80
49	S2	322	C	N3-C2-O2	-5.10	118.33	121.90
4	LA	171	GLY	C-N-CA	5.09	133.00	122.30
1	L5	1963	C	N1-C2-O2	5.09	121.95	118.90
6	LC	2	ALA	C-N-CA	5.08	134.41	121.70
1	L5	469	C	N1-C2-O2	5.08	121.95	118.90
1	L5	4764	A	C6-N1-C2	5.08	121.65	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L5	4766	C	C2-N1-C1'	5.08	124.39	118.80
1	L5	4920	C	N3-C2-O2	-5.08	118.34	121.90
49	S2	1162	C	C6-N1-C2	-5.08	118.27	120.30
1	L5	3838	U	N3-C2-O2	-5.08	118.64	122.20
49	S2	531	A	C4-N9-C1'	5.07	135.42	126.30
49	S2	1309	C	N1-C2-O2	5.07	121.94	118.90
1	L5	1607	C	N1-C2-O2	5.07	121.94	118.90
1	L5	688	U	N3-C2-O2	-5.07	118.65	122.20
1	L5	2892	C	C2-N1-C1'	5.07	124.37	118.80
1	L5	2094	G	C8-N9-C1'	-5.06	120.42	127.00
1	L5	2783	A	C6-N1-C2	5.06	121.64	118.60
1	L5	758	G	C5-C6-O6	5.06	131.64	128.60
1	L5	4561	C	C2-N1-C1'	5.06	124.37	118.80
1	L5	1702	C	N1-C2-O2	5.06	121.94	118.90
49	S2	53	C	C6-N1-C2	-5.06	118.28	120.30
1	L5	4699	U	OP1-P-O3'	5.05	116.32	105.20
49	S2	542	U	N3-C2-O2	-5.05	118.66	122.20
1	L5	2508	U	N3-C2-O2	-5.05	118.66	122.20
49	S2	548	C	C6-N1-C2	-5.05	118.28	120.30
1	L5	2094	G	N3-C4-N9	5.05	129.03	126.00
1	L5	3882	C	C6-N1-C2	-5.05	118.28	120.30
1	L5	4199	C	N1-C2-O2	5.05	121.93	118.90
1	L5	4898	G	C5-C6-O6	5.05	131.63	128.60
49	S2	571	U	N1-C2-O2	5.05	126.33	122.80
1	L5	672	C	C6-N1-C1'	-5.05	114.74	120.80
1	L5	3685	C	C5-C6-N1	5.05	123.52	121.00
1	L5	4112	C	N3-C2-O2	-5.05	118.37	121.90
49	S2	1265	A	N3-C4-C5	-5.05	123.27	126.80
1	L5	4306	U	N3-C2-O2	-5.04	118.67	122.20
49	S2	1292	C	C6-N1-C2	-5.04	118.28	120.30
1	L5	3911	C	C5-C6-N1	5.04	123.52	121.00
1	L5	4215	C	C6-N1-C2	-5.04	118.28	120.30
3	L8	118	C	C2-N1-C1'	5.04	124.34	118.80
49	S2	340	C	N1-C2-O2	5.04	121.92	118.90
1	L5	4766	C	C6-N1-C2	-5.04	118.28	120.30
49	S2	566	U	N3-C2-O2	-5.04	118.67	122.20
49	S2	581	U	C2-N1-C1'	5.04	123.74	117.70
1	L5	4758	U	C6-N1-C1'	-5.03	114.15	121.20
49	S2	402	C	N1-C2-O2	5.03	121.92	118.90
1	L5	3775	A	O4'-C1'-N9	5.03	112.22	108.20
1	L5	4340	U	C5-C6-N1	5.03	125.22	122.70
1	L5	1245	C	C5-C6-N1	5.03	123.51	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
48	Lz	190	LEU	CA-CB-CG	5.02	126.85	115.30
49	S2	1234	C	C2-N1-C1'	5.02	124.33	118.80
84	CC	34	U	C2-N1-C1'	5.02	123.73	117.70
1	L5	704	C	N1-C2-O2	5.02	121.91	118.90
1	L5	2505	C	C5-C6-N1	5.02	123.51	121.00
84	CC	74	C	C5-C6-N1	5.02	123.51	121.00
1	L5	2096	G	N3-C4-N9	5.01	129.01	126.00
1	L5	4940	C	N1-C2-O2	5.01	121.91	118.90
2	L7	76	U	N1-C2-O2	5.01	126.31	122.80
1	L5	1709	C	N1-C2-O2	5.01	121.91	118.90
1	L5	3866	C	C5-C6-N1	5.01	123.51	121.00
1	L5	181	C	C6-N1-C1'	-5.01	114.79	120.80
1	L5	2257	C	N3-C2-O2	-5.01	118.40	121.90
1	L5	489	C	C6-N1-C2	-5.00	118.30	120.30
1	L5	1327	C	C6-N1-C2	-5.00	118.30	120.30
1	L5	4682	U	C2-N1-C1'	5.00	123.70	117.70
49	S2	542	U	C2-N1-C1'	5.00	123.70	117.70

There are no chirality outliers.

All (45) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L5	4906	U	Sidechain
4	LA	110	GLY	Peptide
4	LA	54	ARG	Peptide
5	LB	17	LEU	Peptide
5	LB	258	HIS	Peptide
11	LH	106	GLN	Peptide
13	LJ	94	LEU	Peptide
14	LL	154	VAL	Peptide
15	LM	87	ALA	Peptide
15	LM	88	ALA	Peptide
17	LO	110	PRO	Peptide
18	LP	131	ARG	Peptide
22	LT	135	PRO	Peptide
29	La	22	ILE	Peptide
34	Lf	103	VAL	Peptide
34	Lf	106	TYR	Peptide
36	Lh	86	LYS	Peptide
38	Lj	39	TYR	Peptide
46	Ls	148	SER	Peptide
46	Ls	149	ARG	Peptide

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Mol	Chain	Res	Type	Group
47	Lt	141	CYS	Peptide
47	Lt	147	HIS	Peptide
47	Lt	53	TRP	Peptide
48	Lz	183	ILE	Peptide
51	SB	221	PRO	Peptide
52	SD	164	VAL	Peptide
52	SD	192	TRP	Peptide
52	SD	193	ASP	Peptide
54	SF	126	THR	Peptide
54	SF	78	MET	Peptide
55	SH	15	LYS	Peptide
56	SI	130	THR	Peptide
73	SJ	137	VAL	Peptide
74	SM	43	ASP	Peptide
60	SQ	43	GLU	Peptide
61	SR	84	TYR	Peptide
63	ST	46	ALA	Peptide
65	SV	78	ILE	Peptide
66	SX	125	VAL	Peptide
66	SX	126	ALA	Peptide
66	SX	86	PRO	Peptide
79	SZ	46	ASN	Peptide
79	SZ	78	LYS	Peptide
80	Sb	75	GLU	Peptide
68	Sc	64	GLU	Peptide

## 5.2 Too-close contacts [i](#)

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

## 5.3 Torsion angles [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%)	221 (90%)	24 (10%)	1 (0%)	34	60
5	LB	400/403 (99%)	379 (95%)	19 (5%)	2 (0%)	29	54
6	LC	366/427 (86%)	340 (93%)	25 (7%)	1 (0%)	41	66
7	LD	291/297 (98%)	272 (94%)	19 (6%)	0	100	100
8	LE	232/288 (81%)	210 (90%)	22 (10%)	0	100	100
9	LF	223/248 (90%)	212 (95%)	11 (5%)	0	100	100
10	LG	239/266 (90%)	224 (94%)	15 (6%)	0	100	100
11	LH	188/192 (98%)	173 (92%)	15 (8%)	0	100	100
12	LI	198/214 (92%)	180 (91%)	17 (9%)	1 (0%)	29	54
13	LJ	174/178 (98%)	157 (90%)	17 (10%)	0	100	100
14	LL	208/211 (99%)	190 (91%)	18 (9%)	0	100	100
15	LM	137/215 (64%)	129 (94%)	7 (5%)	1 (1%)	22	46
16	LN	201/204 (98%)	189 (94%)	10 (5%)	2 (1%)	15	37
17	LO	199/203 (98%)	192 (96%)	7 (4%)	0	100	100
18	LP	151/184 (82%)	143 (95%)	7 (5%)	1 (1%)	22	46
19	LQ	185/188 (98%)	177 (96%)	8 (4%)	0	100	100
20	LR	185/196 (94%)	180 (97%)	5 (3%)	0	100	100
21	LS	173/176 (98%)	161 (93%)	12 (7%)	0	100	100
22	LT	157/160 (98%)	147 (94%)	10 (6%)	0	100	100
23	LU	99/128 (77%)	86 (87%)	13 (13%)	0	100	100
24	LV	129/140 (92%)	122 (95%)	7 (5%)	0	100	100
25	LW	122/157 (78%)	114 (93%)	8 (7%)	0	100	100
26	LX	118/156 (76%)	115 (98%)	3 (2%)	0	100	100
27	LY	132/145 (91%)	124 (94%)	8 (6%)	0	100	100
28	LZ	133/136 (98%)	122 (92%)	11 (8%)	0	100	100
29	La	145/148 (98%)	135 (93%)	10 (7%)	0	100	100
30	Lb	105/159 (66%)	97 (92%)	8 (8%)	0	100	100
31	Lc	96/115 (84%)	90 (94%)	6 (6%)	0	100	100
32	Ld	105/125 (84%)	98 (93%)	7 (7%)	0	100	100
33	Le	126/135 (93%)	121 (96%)	4 (3%)	1 (1%)	19	43
34	Lf	107/110 (97%)	98 (92%)	7 (6%)	2 (2%)	8	20
35	Lg	112/117 (96%)	111 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
36	Lh	120/123 (98%)	118 (98%)	2 (2%)	0	100	100
37	Li	100/105 (95%)	95 (95%)	5 (5%)	0	100	100
38	Lj	84/97 (87%)	77 (92%)	6 (7%)	1 (1%)	13	32
39	Lk	67/70 (96%)	62 (92%)	5 (8%)	0	100	100
40	Ll	48/51 (94%)	46 (96%)	2 (4%)	0	100	100
41	Lm	50/128 (39%)	50 (100%)	0	0	100	100
42	Ln	22/25 (88%)	22 (100%)	0	0	100	100
43	Lo	103/106 (97%)	98 (95%)	5 (5%)	0	100	100
44	Lp	89/92 (97%)	82 (92%)	7 (8%)	0	100	100
45	Lr	123/137 (90%)	116 (94%)	7 (6%)	0	100	100
46	Ls	194/317 (61%)	173 (89%)	19 (10%)	2 (1%)	15	37
47	Lt	137/165 (83%)	107 (78%)	27 (20%)	3 (2%)	6	17
48	Lz	215/217 (99%)	165 (77%)	50 (23%)	0	100	100
50	SA	219/295 (74%)	200 (91%)	19 (9%)	0	100	100
51	SB	212/264 (80%)	200 (94%)	12 (6%)	0	100	100
52	SD	225/243 (93%)	202 (90%)	23 (10%)	0	100	100
53	SE	260/263 (99%)	246 (95%)	14 (5%)	0	100	100
54	SF	180/204 (88%)	161 (89%)	19 (11%)	0	100	100
55	SH	182/194 (94%)	159 (87%)	23 (13%)	0	100	100
56	SI	204/208 (98%)	194 (95%)	10 (5%)	0	100	100
57	SK	96/165 (58%)	81 (84%)	15 (16%)	0	100	100
58	SL	151/158 (96%)	137 (91%)	14 (9%)	0	100	100
59	SP	127/145 (88%)	113 (89%)	14 (11%)	0	100	100
60	SQ	142/146 (97%)	126 (89%)	15 (11%)	1 (1%)	22	46
61	SR	133/135 (98%)	121 (91%)	12 (9%)	0	100	100
62	SS	143/152 (94%)	132 (92%)	11 (8%)	0	100	100
63	ST	141/145 (97%)	127 (90%)	13 (9%)	1 (1%)	22	46
64	SU	102/119 (86%)	93 (91%)	9 (9%)	0	100	100
65	SV	81/83 (98%)	74 (91%)	6 (7%)	1 (1%)	13	32
66	SX	139/143 (97%)	129 (93%)	7 (5%)	3 (2%)	6	17
67	Sa	100/115 (87%)	92 (92%)	7 (7%)	1 (1%)	15	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
68	Sc	62/69 (90%)	51 (82%)	11 (18%)	0	100	100
69	Sd	53/56 (95%)	50 (94%)	2 (4%)	1 (2%)	8	20
70	Sg	311/317 (98%)	267 (86%)	44 (14%)	0	100	100
71	SC	220/293 (75%)	206 (94%)	14 (6%)	0	100	100
72	SG	235/249 (94%)	221 (94%)	14 (6%)	0	100	100
73	SJ	183/194 (94%)	171 (93%)	11 (6%)	1 (0%)	29	54
74	SM	120/132 (91%)	112 (93%)	8 (7%)	0	100	100
75	SN	148/151 (98%)	144 (97%)	4 (3%)	0	100	100
76	SO	138/151 (91%)	124 (90%)	14 (10%)	0	100	100
77	SW	127/130 (98%)	121 (95%)	6 (5%)	0	100	100
78	SY	129/133 (97%)	124 (96%)	5 (4%)	0	100	100
79	SZ	73/125 (58%)	61 (84%)	10 (14%)	2 (3%)	5	12
80	Sb	81/84 (96%)	73 (90%)	8 (10%)	0	100	100
81	Se	56/59 (95%)	52 (93%)	4 (7%)	0	100	100
82	Sf	65/156 (42%)	55 (85%)	10 (15%)	0	100	100
83	CA	350/394 (89%)	326 (93%)	24 (7%)	0	100	100
85	CE	120/223 (54%)	114 (95%)	6 (5%)	0	100	100
86	CF	27/180 (15%)	25 (93%)	2 (7%)	0	100	100
All	All	12369/14184 (87%)	11404 (92%)	936 (8%)	29 (0%)	50	73

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
16	LN	124	ASP
46	Ls	150	GLY
66	SX	127	ASN
15	LM	88	ALA
47	Lt	148	PRO
63	ST	41	LYS
67	Sa	47	ALA
79	SZ	45	ASN
5	LB	302	ASN
34	Lf	80	ASN
66	SX	126	ALA
47	Lt	147	HIS
60	SQ	44	PRO

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Mol	Chain	Res	Type
5	LB	360	LEU
33	Le	73	GLY
34	Lf	107	PRO
47	Lt	121	LEU
79	SZ	78	LYS
4	LA	55	GLY
16	LN	83	LYS
69	Sd	14	PHE
46	Ls	186	GLY
66	SX	78	GLY
38	Lj	40	PRO
73	SJ	123	ILE
6	LC	232	VAL
18	LP	77	GLY
65	SV	79	VAL
12	LI	15	LYS

### 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%)	188 (99%)	2 (1%)	73	90
5	LB	348/349 (100%)	347 (100%)	1 (0%)	92	98
6	LC	306/348 (88%)	304 (99%)	2 (1%)	84	94
7	LD	246/250 (98%)	246 (100%)	0	100	100
8	LE	209/252 (83%)	208 (100%)	1 (0%)	88	96
9	LF	194/215 (90%)	194 (100%)	0	100	100
10	LG	203/223 (91%)	201 (99%)	2 (1%)	76	91
11	LH	169/171 (99%)	168 (99%)	1 (1%)	86	95
12	LI	172/181 (95%)	172 (100%)	0	100	100
13	LJ	148/149 (99%)	146 (99%)	2 (1%)	67	86
14	LL	176/177 (99%)	173 (98%)	3 (2%)	60	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
15	LM	118/161 (73%)	117 (99%)	1 (1%)	81	93
16	LN	171/172 (99%)	171 (100%)	0	100	100
17	LO	173/174 (99%)	172 (99%)	1 (1%)	86	95
18	LP	134/163 (82%)	134 (100%)	0	100	100
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%)	155 (99%)	1 (1%)	86	95
22	LT	139/140 (99%)	138 (99%)	1 (1%)	84	94
23	LU	91/115 (79%)	90 (99%)	1 (1%)	73	90
24	LV	101/107 (94%)	100 (99%)	1 (1%)	76	91
25	LW	103/126 (82%)	101 (98%)	2 (2%)	57	82
26	LX	108/133 (81%)	108 (100%)	0	100	100
27	LY	124/135 (92%)	121 (98%)	3 (2%)	49	77
28	LZ	117/118 (99%)	117 (100%)	0	100	100
29	La	120/121 (99%)	119 (99%)	1 (1%)	81	93
30	Lb	88/126 (70%)	88 (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%)	97 (99%)	1 (1%)	76	91
36	Lh	109/110 (99%)	109 (100%)	0	100	100
37	Li	86/89 (97%)	85 (99%)	1 (1%)	71	88
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%)	47 (100%)	0	100	100
41	Lm	48/116 (41%)	48 (100%)	0	100	100
42	Ln	23/24 (96%)	23 (100%)	0	100	100
43	Lo	93/94 (99%)	93 (100%)	0	100	100
44	Lp	74/75 (99%)	74 (100%)	0	100	100
45	Lr	109/121 (90%)	109 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
46	Ls	162/258 (63%)	160 (99%)	2 (1%)	71	88
47	Lt	112/137 (82%)	111 (99%)	1 (1%)	78	92
48	Lz	195/196 (100%)	187 (96%)	8 (4%)	30	59
50	SA	183/243 (75%)	182 (100%)	1 (0%)	88	96
51	SB	195/231 (84%)	194 (100%)	1 (0%)	88	96
52	SD	190/202 (94%)	188 (99%)	2 (1%)	73	90
53	SE	224/225 (100%)	224 (100%)	0	100	100
54	SF	156/170 (92%)	155 (99%)	1 (1%)	86	95
55	SH	166/174 (95%)	164 (99%)	2 (1%)	71	88
56	SI	178/180 (99%)	176 (99%)	2 (1%)	73	90
57	SK	89/136 (65%)	88 (99%)	1 (1%)	73	90
58	SL	137/142 (96%)	136 (99%)	1 (1%)	84	94
59	SP	115/130 (88%)	114 (99%)	1 (1%)	78	92
60	SQ	119/121 (98%)	118 (99%)	1 (1%)	81	93
61	SR	122/122 (100%)	122 (100%)	0	100	100
62	SS	126/132 (96%)	125 (99%)	1 (1%)	81	93
63	ST	113/115 (98%)	111 (98%)	2 (2%)	59	83
64	SU	94/107 (88%)	94 (100%)	0	100	100
65	SV	67/67 (100%)	66 (98%)	1 (2%)	65	86
66	SX	113/115 (98%)	112 (99%)	1 (1%)	78	92
67	Sa	89/98 (91%)	89 (100%)	0	100	100
68	Sc	57/62 (92%)	57 (100%)	0	100	100
69	Sd	48/49 (98%)	47 (98%)	1 (2%)	53	80
70	Sg	272/275 (99%)	267 (98%)	5 (2%)	59	83
71	SC	188/225 (84%)	186 (99%)	2 (1%)	73	90
72	SG	207/218 (95%)	203 (98%)	4 (2%)	57	82
73	SJ	161/168 (96%)	161 (100%)	0	100	100
74	SM	102/108 (94%)	98 (96%)	4 (4%)	32	61
75	SN	130/131 (99%)	130 (100%)	0	100	100
76	SO	110/119 (92%)	109 (99%)	1 (1%)	78	92
77	SW	112/113 (99%)	112 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
78	SY	113/115 (98%)	109 (96%)	4 (4%)	36	65
79	SZ	66/103 (64%)	63 (96%)	3 (4%)	27	55
80	Sb	75/76 (99%)	75 (100%)	0	100	100
81	Se	42/48 (88%)	41 (98%)	1 (2%)	49	77
82	Sf	60/140 (43%)	59 (98%)	1 (2%)	60	84
83	CA	303/336 (90%)	302 (100%)	1 (0%)	92	98
85	CE	108/190 (57%)	104 (96%)	4 (4%)	34	63
86	CF	26/151 (17%)	26 (100%)	0	100	100
All	All	10766/12069 (89%)	10678 (99%)	88 (1%)	82	93

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	LA	30	ARG
4	LA	207	VAL
5	LB	258	HIS
6	LC	257	PHE
6	LC	364	LYS
8	LE	56	ARG
10	LG	111	LYS
10	LG	175	ARG
11	LH	17	ASP
13	LJ	95	ARG
13	LJ	118	LYS
14	LL	59	VAL
14	LL	67	HIS
14	LL	144	LEU
15	LM	38	VAL
17	LO	117	ARG
21	LS	90	THR
22	LT	36	LYS
23	LU	67	LYS
24	LV	48	ARG
25	LW	77	LYS
25	LW	100	VAL
27	LY	74	TYR
27	LY	79	VAL
27	LY	84	ARG
29	La	92	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
35	Lg	54	ARG
37	Li	29	ARG
46	Ls	77	LYS
46	Ls	149	ARG
47	Lt	16	ARG
48	Lz	7	ARG
48	Lz	17	VAL
48	Lz	47	LYS
48	Lz	85	MET
48	Lz	102	LYS
48	Lz	139	THR
48	Lz	145	VAL
48	Lz	215	ARG
50	SA	73	ASP
51	SB	62	LEU
52	SD	76	ARG
52	SD	167	TYR
54	SF	45	TYR
55	SH	32	MET
55	SH	69	LEU
56	SI	47	ARG
56	SI	125	LYS
57	SK	98	ARG
58	SL	69	ARG
59	SP	70	MET
60	SQ	71	ARG
62	SS	142	ARG
63	ST	24	LYS
63	ST	41	LYS
65	SV	83	PHE
66	SX	105	PHE
69	Sd	27	ARG
70	Sg	94	THR
70	Sg	185	LYS
70	Sg	225	LYS
70	Sg	248	LEU
70	Sg	289	LEU
71	SC	64	THR
71	SC	137	VAL
72	SG	22	ARG
72	SG	98	ARG
72	SG	119	LYS

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Mol	Chain	Res	Type
72	SG	215	LYS
74	SM	18	LEU
74	SM	46	GLN
74	SM	63	LYS
74	SM	66	GLU
76	SO	150	ARG
78	SY	16	ARG
78	SY	26	ASP
78	SY	46	LYS
78	SY	99	LYS
79	SZ	41	ARG
79	SZ	50	PHE
79	SZ	92	LEU
81	Se	26	LYS
82	Sf	104	LYS
83	CA	51	LYS
85	CE	29	LYS
85	CE	30	LYS
85	CE	33	GLU
85	CE	199	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (104) such sidechains are listed below:

Mol	Chain	Res	Type
4	LA	97	ASN
4	LA	132	ASN
5	LB	175	GLN
5	LB	213	GLN
6	LC	21	ASN
6	LC	38	ASN
6	LC	178	ASN
7	LD	282	GLN
7	LD	291	GLN
8	LE	190	HIS
9	LF	63	GLN
9	LF	206	ASN
10	LG	94	GLN
10	LG	112	GLN
15	LM	125	ASN
16	LN	29	GLN
16	LN	109	HIS
16	LN	149	GLN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
18	LP	21	ASN
18	LP	34	GLN
18	LP	80	GLN
18	LP	97	ASN
18	LP	116	HIS
19	LQ	21	GLN
19	LQ	44	ASN
19	LQ	125	GLN
19	LQ	188	ASN
20	LR	39	GLN
20	LR	130	ASN
21	LS	77	ASN
23	LU	17	GLN
23	LU	50	ASN
25	LW	48	GLN
25	LW	96	GLN
25	LW	120	GLN
26	LX	151	ASN
27	LY	40	GLN
28	LZ	78	ASN
29	La	67	GLN
33	Le	107	ASN
34	Lf	56	ASN
36	Lh	101	ASN
37	Li	12	ASN
38	Lj	66	HIS
43	Lo	19	GLN
43	Lo	45	GLN
43	Lo	51	GLN
44	Lp	56	HIS
45	Lr	70	GLN
45	Lr	100	ASN
46	Ls	127	ASN
46	Ls	176	ASN
47	Lt	70	GLN
48	Lz	35	GLN
48	Lz	182	ASN
50	SA	141	ASN
51	SB	179	ASN
51	SB	202	GLN
53	SE	98	ASN
53	SE	112	HIS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
53	SE	138	HIS
54	SF	148	ASN
55	SH	12	ASN
55	SH	33	ASN
55	SH	114	GLN
56	SI	9	HIS
56	SI	84	ASN
57	SK	7	ASN
57	SK	42	ASN
57	SK	44	HIS
57	SK	50	GLN
58	SL	65	ASN
58	SL	83	GLN
59	SP	103	ASN
59	SP	137	HIS
61	SR	31	ASN
65	SV	49	GLN
65	SV	82	ASN
67	Sa	8	ASN
69	Sd	5	GLN
69	Sd	26	ASN
70	Sg	4	GLN
70	Sg	14	HIS
70	Sg	117	ASN
71	SC	267	GLN
72	SG	110	ASN
72	SG	146	ASN
73	SJ	125	HIS
75	SN	49	GLN
77	SW	90	GLN
77	SW	91	ASN
80	Sb	26	GLN
80	Sb	65	GLN
81	Se	3	HIS
81	Se	58	ASN
83	CA	83	ASN
83	CA	134	GLN
83	CA	203	GLN
83	CA	204	ASN
83	CA	254	GLN
85	CE	51	GLN
85	CE	63	GLN

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Mol	Chain	Res	Type
85	CE	71	GLN
85	CE	180	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	L5	3706/5066 (73%)	856 (23%)	20 (0%)
2	L7	119/121 (98%)	13 (10%)	0
3	L8	155/157 (98%)	28 (18%)	0
49	S2	1717/1869 (91%)	382 (22%)	7 (0%)
84	CC	74/75 (98%)	23 (31%)	2 (2%)
All	All	5771/7288 (79%)	1302 (22%)	29 (0%)

All (1302) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	L5	2	G
1	L5	17	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	69	A
1	L5	72	C
1	L5	73	A
1	L5	74	G
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	119	G
1	L5	120	A
1	L5	132	G
1	L5	133	C
1	L5	134	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	135	G
1	L5	136	C
1	L5	137	G
1	L5	152	U
1	L5	159	C
1	L5	165	A
1	L5	183	C
1	L5	184	U
1	L5	185	C
1	L5	187	U
1	L5	188	G
1	L5	189	G
1	L5	200	U
1	L5	207	G
1	L5	210	C
1	L5	218	A
1	L5	219	G
1	L5	233	U
1	L5	234	G
1	L5	255	C
1	L5	256	G
1	L5	258	G
1	L5	261	G
1	L5	265	C
1	L5	266	C
1	L5	267	G
1	L5	269	G
1	L5	278	G
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	349	A
1	L5	360	A
1	L5	373	G
1	L5	387	G
1	L5	388	A
1	L5	398	A
1	L5	399	G
1	L5	407	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
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	449	C
1	L5	450	G
1	L5	452	A
1	L5	453	G
1	L5	454	U
1	L5	456	C
1	L5	457	G
1	L5	465	G
1	L5	467	U
1	L5	468	U
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	499	G
1	L5	500	G
1	L5	502	C
1	L5	503	C
1	L5	504	G
1	L5	505	G
1	L5	507	G
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	518	G
1	L5	643	C

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	646	G
1	L5	656	C
1	L5	657	C
1	L5	658	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	673	C
1	L5	685	C
1	L5	686	A
1	L5	687	U
1	L5	688	U
1	L5	696	C
1	L5	703	G
1	L5	704	C
1	L5	708	G
1	L5	731	G
1	L5	738	C
1	L5	739	G
1	L5	742	G
1	L5	746	A
1	L5	753	C
1	L5	758	G
1	L5	759	G
1	L5	760	G
1	L5	904	C
1	L5	913	U
1	L5	914	U
1	L5	915	A
1	L5	916	C
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

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	941	C
1	L5	945	U
1	L5	958	G
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	977	C
1	L5	982	U
1	L5	985	C
1	L5	989	U
1	L5	990	C
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	1074	G
1	L5	1075	G
1	L5	1082	C
1	L5	1083	U
1	L5	1095	A
1	L5	1168	G
1	L5	1170	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

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
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	1218	G
1	L5	1219	G
1	L5	1222	A
1	L5	1241	C
1	L5	1245	C
1	L5	1246	G
1	L5	1253	G
1	L5	1254	A
1	L5	1257	A
1	L5	1258	G
1	L5	1266	G
1	L5	1267	C
1	L5	1269	G
1	L5	1270	A
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	1301	C
1	L5	1302	U
1	L5	1304	C
1	L5	1312	A
1	L5	1324	A
1	L5	1326	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	1337	A
1	L5	1344	C
1	L5	1354	A
1	L5	1358	G
1	L5	1359	G
1	L5	1364	U
1	L5	1365	C
1	L5	1367	C
1	L5	1378	C
1	L5	1379	C
1	L5	1387	A
1	L5	1393	G
1	L5	1394	G
1	L5	1397	A
1	L5	1403	G
1	L5	1404	G
1	L5	1405	C
1	L5	1407	C
1	L5	1409	C
1	L5	1410	U
1	L5	1411	C
1	L5	1414	C
1	L5	1415	G
1	L5	1417	C
1	L5	1420	A
1	L5	1435	G
1	L5	1437	C
1	L5	1439	C
1	L5	1440	U
1	L5	1442	C
1	L5	1445	U
1	L5	1446	C
1	L5	1447	C
1	L5	1482	G
1	L5	1483	C
1	L5	1493	G
1	L5	1497	A
1	L5	1498	G
1	L5	1502	G
1	L5	1517	G
1	L5	1518	A
1	L5	1525	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	1534	A
1	L5	1547	A
1	L5	1562	G
1	L5	1566	C
1	L5	1578	U
1	L5	1591	U
1	L5	1596	U
1	L5	1613	A
1	L5	1624	G
1	L5	1625	G
1	L5	1631	A
1	L5	1633	G
1	L5	1634	A
1	L5	1640	C
1	L5	1641	G
1	L5	1642	A
1	L5	1654	G
1	L5	1661	C
1	L5	1676	C
1	L5	1677	U
1	L5	1678	C
1	L5	1697	G
1	L5	1698	C
1	L5	1699	A
1	L5	1700	G
1	L5	1703	C
1	L5	1704	C
1	L5	1705	G
1	L5	1707	C
1	L5	1709	C
1	L5	1715	C
1	L5	1716	G
1	L5	1719	A
1	L5	1726	U
1	L5	1734	G
1	L5	1735	U
1	L5	1741	G
1	L5	1742	A
1	L5	1755	C
1	L5	1757	U
1	L5	1758	G
1	L5	1760	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
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	1772	C
1	L5	1775	A
1	L5	1785	C
1	L5	1787	A
1	L5	1792	U
1	L5	1797	G
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
1	L5	1892	A
1	L5	1897	A
1	L5	1898	C
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	1931	C
1	L5	1932	A
1	L5	1936	C
1	L5	1940	G
1	L5	1948	G
1	L5	1951	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	1959	U
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	2003	G
1	L5	2004	U
1	L5	2011	C
1	L5	2014	C
1	L5	2017	A
1	L5	2018	C
1	L5	2024	G
1	L5	2025	A
1	L5	2026	A
1	L5	2033	A
1	L5	2034	G
1	L5	2042	A
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

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	2095	A
1	L5	2096	G
1	L5	2097	U
1	L5	2098	G
1	L5	2101	C
1	L5	2102	G
1	L5	2110	C
1	L5	2112	G
1	L5	2113	G
1	L5	2250	C
1	L5	2252	G
1	L5	2253	A
1	L5	2256	C
1	L5	2258	C
1	L5	2260	C
1	L5	2263	A
1	L5	2289	C
1	L5	2300	A
1	L5	2301	G
1	L5	2306	G
1	L5	2313	A
1	L5	2332	A
1	L5	2333	G
1	L5	2348	G
1	L5	2351	C
1	L5	2360	A
1	L5	2369	U
1	L5	2383	C
1	L5	2395	A
1	L5	2397	G
1	L5	2417	A
1	L5	2418	A
1	L5	2421	G
1	L5	2425	U
1	L5	2437	C
1	L5	2441	C
1	L5	2450	G
1	L5	2453	A
1	L5	2464	C
1	L5	2465	C
1	L5	2469	C
1	L5	2474	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	2475	G
1	L5	2478	C
1	L5	2479	G
1	L5	2483	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	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	2537	A
1	L5	2544	G
1	L5	2546	G
1	L5	2547	G
1	L5	2554	U
1	L5	2559	G
1	L5	2560	C
1	L5	2573	A
1	L5	2583	C
1	L5	2586	G
1	L5	2587	A
1	L5	2589	C
1	L5	2601	A
1	L5	2611	A
1	L5	2618	G
1	L5	2627	C
1	L5	2652	G
1	L5	2653	C
1	L5	2662	G
1	L5	2669	C
1	L5	2670	C
1	L5	2675	G
1	L5	2676	A
1	L5	2687	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
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	2725	A
1	L5	2726	G
1	L5	2739	C
1	L5	2742	G
1	L5	2743	A
1	L5	2746	A
1	L5	2756	G
1	L5	2761	U
1	L5	2763	U
1	L5	2764	A
1	L5	2769	U
1	L5	2770	C
1	L5	2787	A
1	L5	2788	U
1	L5	2790	U
1	L5	2799	G
1	L5	2814	C
1	L5	2815	A
1	L5	2826	U
1	L5	2827	G
1	L5	2835	A
1	L5	2838	G
1	L5	2842	G
1	L5	2848	G
1	L5	2855	G
1	L5	2867	C
1	L5	2877	G
1	L5	2892	C
1	L5	2900	U
1	L5	2902	G
1	L5	2903	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	2904	U
1	L5	2905	C
1	L5	2906	G
1	L5	2907	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	3605	C
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	3662	A
1	L5	3664	G
1	L5	3670	C
1	L5	3673	C
1	L5	3674	G
1	L5	3711	A
1	L5	3713	U
1	L5	3714	G
1	L5	3727	A
1	L5	3729	U
1	L5	3735	G
1	L5	3736	A
1	L5	3748	A
1	L5	3750	G
1	L5	3753	G
1	L5	3757	G
1	L5	3758	U
1	L5	3759	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	3760	A
1	L5	3761	C
1	L5	3771	C
1	L5	3776	G
1	L5	3777	G
1	L5	3783	A
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	3930	U
1	L5	3938	G
1	L5	3939	G
1	L5	3942	A
1	L5	3943	A
1	L5	3944	G
1	L5	3947	A
1	L5	3950	U
1	L5	3953	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	3955	G
1	L5	3956	G
1	L5	3957	U
1	L5	3958	G
1	L5	3959	U
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	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	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	4056	A
1	L5	4057	C
1	L5	4058	U
1	L5	4059	C
1	L5	4062	A
1	L5	4063	U
1	L5	4064	C
1	L5	4065	G
1	L5	4076	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	4084	G
1	L5	4086	G
1	L5	4091	G
1	L5	4096	C
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	4122	G
1	L5	4127	A
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	4160	C
1	L5	4162	C
1	L5	4163	U
1	L5	4168	G
1	L5	4170	A
1	L5	4177	C
1	L5	4183	G
1	L5	4184	G
1	L5	4191	G
1	L5	4196	G
1	L5	4197	G
1	L5	4203	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	4222	G
1	L5	4225	G
1	L5	4228	G
1	L5	4229	U
1	L5	4233	A
1	L5	4249	G
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	4295	U
1	L5	4304	A
1	L5	4305	G
1	L5	4306	U
1	L5	4314	C
1	L5	4319	C
1	L5	4329	G
1	L5	4330	G
1	L5	4332	C
1	L5	4339	A
1	L5	4349	C
1	L5	4371	G
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	4386	C
1	L5	4387	C
1	L5	4391	G
1	L5	4393	G
1	L5	4394	A
1	L5	4398	C
1	L5	4422	A
1	L5	4426	C
1	L5	4448	G
1	L5	4449	A
1	L5	4452	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
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	4515	G
1	L5	4519	C
1	L5	4524	G
1	L5	4531	U
1	L5	4545	G
1	L5	4548	A
1	L5	4549	G
1	L5	4557	U
1	L5	4560	C
1	L5	4567	G
1	L5	4573	G
1	L5	4575	G
1	L5	4589	A
1	L5	4590	A
1	L5	4600	G
1	L5	4617	G
1	L5	4635	A
1	L5	4636	U
1	L5	4637	G
1	L5	4652	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	4693	C
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

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	4730	C
1	L5	4731	G
1	L5	4734	A
1	L5	4740	G
1	L5	4741	C
1	L5	4742	G
1	L5	4745	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	4771	C
1	L5	4772	C
1	L5	4775	C
1	L5	4859	C
1	L5	4870	G
1	L5	4871	C
1	L5	4875	G
1	L5	4880	C
1	L5	4881	U
1	L5	4882	U
1	L5	4883	C
1	L5	4888	U
1	L5	4889	G
1	L5	4895	C
1	L5	4896	G
1	L5	4900	C
1	L5	4901	G
1	L5	4902	C
1	L5	4904	G
1	L5	4905	C
1	L5	4906	U
1	L5	4910	A
1	L5	4912	G
1	L5	4914	C
1	L5	4922	C
1	L5	4923	C
1	L5	4924	C
1	L5	4925	U
1	L5	4926	C

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	4927	G
1	L5	4928	C
1	L5	4934	A
1	L5	4937	C
1	L5	4940	C
1	L5	4941	G
1	L5	4943	A
1	L5	4951	G
1	L5	4960	G
1	L5	4961	G
1	L5	4966	A
1	L5	4976	U
1	L5	4985	U
1	L5	4988	U
1	L5	4989	U
1	L5	4991	U
1	L5	5014	A
1	L5	5017	G
1	L5	5022	U
1	L5	5025	C
1	L5	5026	U
1	L5	5027	C
1	L5	5028	G
1	L5	5029	C
1	L5	5030	U
1	L5	5034	A
1	L5	5041	G
1	L5	5047	C
1	L5	5050	C
1	L5	5054	C
1	L5	5058	A
1	L5	5061	A
1	L5	5069	U
2	L7	4	U
2	L7	5	A
2	L7	33	U
2	L7	38	U
2	L7	49	A
2	L7	53	U
2	L7	54	A
2	L7	63	C
2	L7	64	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
2	L7	97	G
2	L7	100	A
2	L7	110	G
2	L7	111	C
3	L8	25	G
3	L8	34	U
3	L8	35	C
3	L8	48	A
3	L8	59	A
3	L8	60	G
3	L8	62	A
3	L8	63	U
3	L8	80	A
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
3	L8	110	U
3	L8	111	U
3	L8	114	G
3	L8	123	U
3	L8	124	U
3	L8	125	C
3	L8	126	C
3	L8	127	U
3	L8	153	C
3	L8	156	U
49	S2	13	C
49	S2	17	C
49	S2	33	G
49	S2	41	G
49	S2	44	U
49	S2	45	A
49	S2	46	A
49	S2	56	G
49	S2	58	C
49	S2	59	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	64	A
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	121	U
49	S2	126	G
49	S2	130	G
49	S2	139	C
49	S2	143	U
49	S2	147	A
49	S2	149	A
49	S2	155	G
49	S2	158	A
49	S2	159	A
49	S2	160	U
49	S2	161	U
49	S2	162	C
49	S2	175	A
49	S2	179	C
49	S2	182	C
49	S2	184	G
49	S2	190	G
49	S2	196	C
49	S2	197	U
49	S2	198	U
49	S2	199	C
49	S2	200	G
49	S2	203	G
49	S2	204	G
49	S2	206	G
49	S2	208	G
49	S2	209	A
49	S2	214	U
49	S2	290	U
49	S2	291	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	292	A
49	S2	295	C
49	S2	302	A
49	S2	306	C
49	S2	307	G
49	S2	308	G
49	S2	309	G
49	S2	311	C
49	S2	318	A
49	S2	319	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	332	G
49	S2	338	G
49	S2	339	A
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	407	G
49	S2	408	A
49	S2	409	C
49	S2	417	C
49	S2	418	A
49	S2	429	C
49	S2	438	G
49	S2	448	A
49	S2	449	A
49	S2	450	C
49	S2	452	G
49	S2	464	A
49	S2	465	A
49	S2	466	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	471	G
49	S2	472	C
49	S2	473	A
49	S2	474	G
49	S2	482	G
49	S2	483	C
49	S2	487	U
49	S2	488	U
49	S2	492	C
49	S2	493	A
49	S2	500	A
49	S2	502	C
49	S2	512	A
49	S2	516	A
49	S2	525	A
49	S2	529	A
49	S2	531	A
49	S2	532	C
49	S2	533	A
49	S2	536	A
49	S2	537	C
49	S2	540	U
49	S2	542	U
49	S2	546	G
49	S2	547	G
49	S2	556	U
49	S2	557	U
49	S2	558	G
49	S2	559	G
49	S2	563	G
49	S2	564	A
49	S2	576	A
49	S2	583	A
49	S2	584	A
49	S2	587	A
49	S2	589	G
49	S2	590	A
49	S2	591	U
49	S2	597	G
49	S2	604	A
49	S2	607	U
49	S2	608	C

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	614	C
49	S2	623	G
49	S2	628	A
49	S2	629	A
49	S2	631	U
49	S2	643	A
49	S2	644	G
49	S2	655	A
49	S2	660	C
49	S2	668	A
49	S2	669	A
49	S2	671	A
49	S2	672	A
49	S2	673	G
49	S2	683	G
49	S2	687	C
49	S2	688	U
49	S2	689	U
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	734	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	G
49	S2	810	A
49	S2	811	A
49	S2	821	G
49	S2	822	U
49	S2	823	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	824	C
49	S2	830	A
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	842	C
49	S2	847	A
49	S2	869	A
49	S2	870	A
49	S2	873	G
49	S2	874	G
49	S2	878	G
49	S2	880	G
49	S2	882	U
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	894	G
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	913	A
49	S2	914	U
49	S2	920	A
49	S2	930	C
49	S2	933	G
49	S2	934	G
49	S2	943	U
49	S2	963	A
49	S2	971	G
49	S2	972	A
49	S2	990	A

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	992	A
49	S2	999	G
49	S2	1001	A
49	S2	1017	U
49	S2	1018	U
49	S2	1023	A
49	S2	1027	A
49	S2	1028	A
49	S2	1033	G
49	S2	1061	U
49	S2	1062	A
49	S2	1083	A
49	S2	1085	C
49	S2	1089	G
49	S2	1109	C
49	S2	1114	U
49	S2	1115	U
49	S2	1116	C
49	S2	1118	C
49	S2	1121	G
49	S2	1138	C
49	S2	1153	C
49	S2	1154	U
49	S2	1195	A
49	S2	1207	G
49	S2	1208	A
49	S2	1215	C
49	S2	1216	C
49	S2	1217	A
49	S2	1224	G
49	S2	1227	G
49	S2	1242	U
49	S2	1243	U
49	S2	1251	A
49	S2	1253	A
49	S2	1256	G
49	S2	1257	G
49	S2	1259	A
49	S2	1264	C
49	S2	1265	A
49	S2	1274	G
49	S2	1275	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	1283	C
49	S2	1284	A
49	S2	1286	G
49	S2	1294	G
49	S2	1295	A
49	S2	1302	G
49	S2	1306	U
49	S2	1308	U
49	S2	1312	G
49	S2	1342	U
49	S2	1343	U
49	S2	1348	G
49	S2	1371	U
49	S2	1372	U
49	S2	1373	C
49	S2	1378	A
49	S2	1402	A
49	S2	1404	U
49	S2	1415	C
49	S2	1419	C
49	S2	1420	G
49	S2	1421	A
49	S2	1422	G
49	S2	1423	C
49	S2	1424	G
49	S2	1433	C
49	S2	1435	C
49	S2	1436	C
49	S2	1438	A
49	S2	1452	A
49	S2	1454	A
49	S2	1455	A
49	S2	1462	U
49	S2	1463	U
49	S2	1480	A
49	S2	1487	A
49	S2	1489	A
49	S2	1490	G
49	S2	1495	G
49	S2	1497	G
49	S2	1498	A
49	S2	1505	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	1507	G
49	S2	1509	U
49	S2	1520	G
49	S2	1521	C
49	S2	1533	A
49	S2	1537	A
49	S2	1544	C
49	S2	1552	G
49	S2	1553	C
49	S2	1556	A
49	S2	1558	C
49	S2	1570	G
49	S2	1574	C
49	S2	1578	U
49	S2	1580	A
49	S2	1585	U
49	S2	1587	G
49	S2	1588	A
49	S2	1594	A
49	S2	1601	A
49	S2	1621	U
49	S2	1623	A
49	S2	1634	A
49	S2	1638	G
49	S2	1644	C
49	S2	1647	A
49	S2	1648	G
49	S2	1654	G
49	S2	1661	A
49	S2	1663	A
49	S2	1664	A
49	S2	1665	G
49	S2	1671	G
49	S2	1680	G
49	S2	1683	C
49	S2	1686	G
49	S2	1698	C
49	S2	1699	A
49	S2	1719	A
49	S2	1721	U
49	S2	1722	G
49	S2	1743	G

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
49	S2	1744	G
49	S2	1745	A
49	S2	1752	C
49	S2	1753	C
49	S2	1754	G
49	S2	1757	G
49	S2	1758	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	1798	C
49	S2	1822	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
84	CC	11	C
84	CC	16	C
84	CC	17	G
84	CC	18	G
84	CC	19	C
84	CC	20	U

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
84	CC	21	A
84	CC	23	C
84	CC	28	C
84	CC	32	C
84	CC	33	U
84	CC	35	U
84	CC	36	C
84	CC	37	A
84	CC	39	C
84	CC	42	G
84	CC	46	A
84	CC	47	C
84	CC	48	C
84	CC	55	C
84	CC	58	C
84	CC	72	G
84	CC	75	A

All (29) RNA pucker outliers are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	L5	406	C
1	L5	914	U
1	L5	955	G
1	L5	1082	C
1	L5	1633	G
1	L5	1733	G
1	L5	1977	C
1	L5	2033	A
1	L5	2112	G
1	L5	2416	G
1	L5	2675	G
1	L5	2760	G
1	L5	2786	C
1	L5	3614	G
1	L5	3673	C
1	L5	3876	A
1	L5	4378	A
1	L5	4699	U
1	L5	4904	G
1	L5	4913	G
49	S2	112	U

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Mol	Chain	Res	Type
49	S2	291	G
49	S2	417	C
49	S2	563	G
49	S2	688	U
49	S2	1265	A
49	S2	1434	C
84	CC	35	U
84	CC	74	C

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

There are no non-standard protein/DNA/RNA residues in this entry.

#### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

#### 5.6 Ligand geometry [i](#)

Of 264 ligands modelled in this entry, 264 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

#### 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

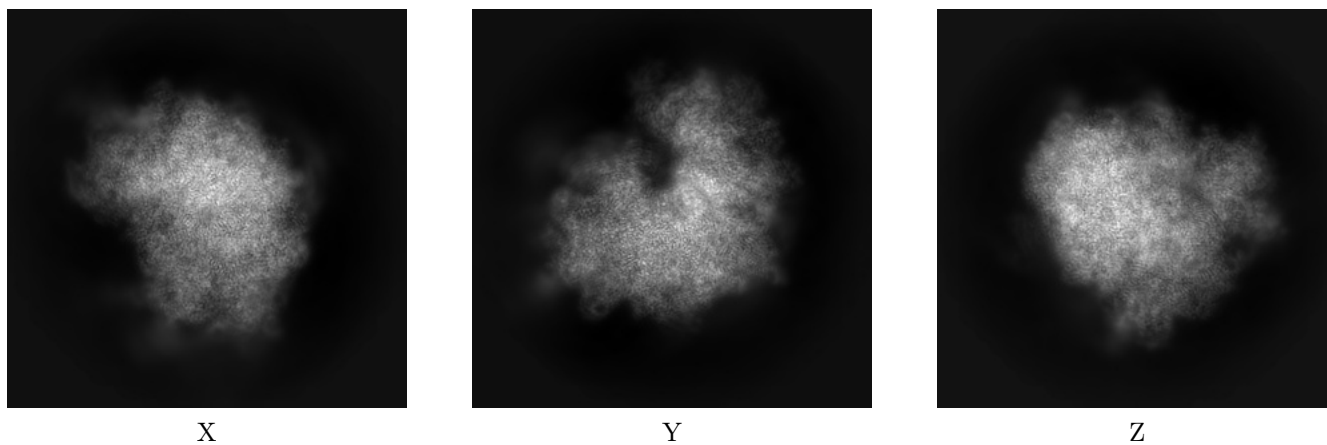
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11288. These allow visual inspection of the internal detail of the map and identification of artifacts.

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

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

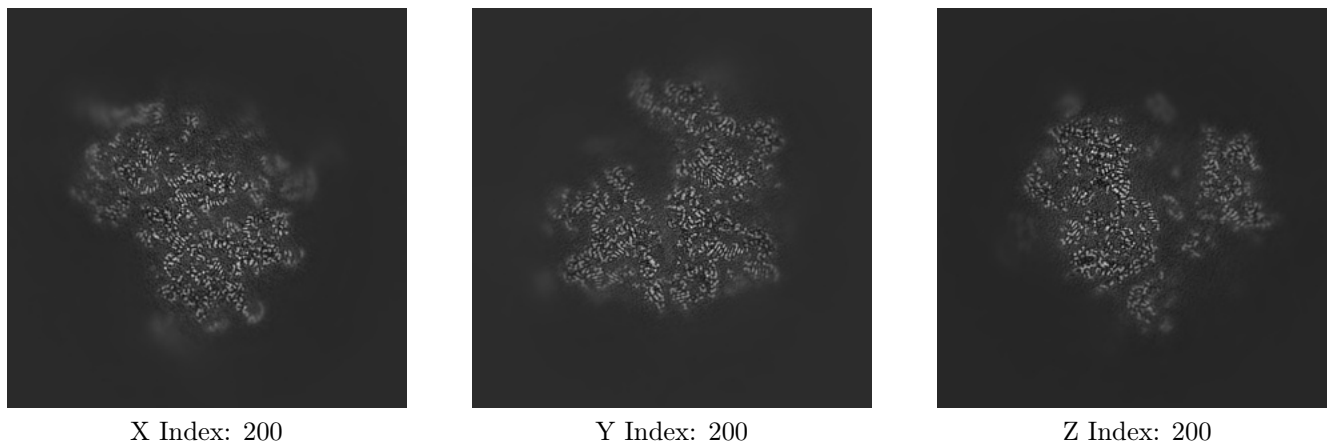
#### 6.1.1 Primary map



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

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

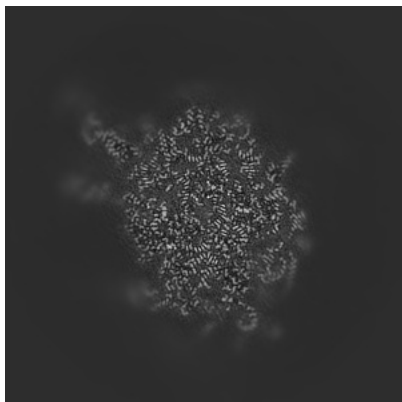
#### 6.2.1 Primary map



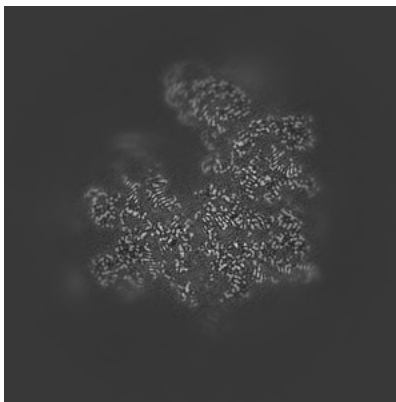
The images above show central slices of the map in three orthogonal directions.

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

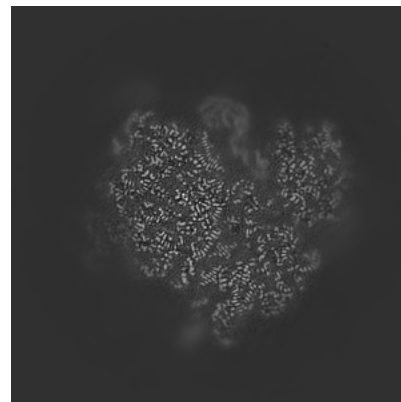
### 6.3.1 Primary map



X Index: 178



Y Index: 194



Z Index: 224

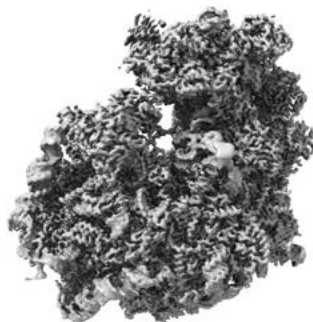
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

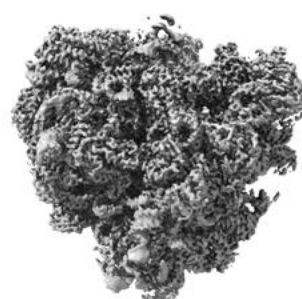
### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.



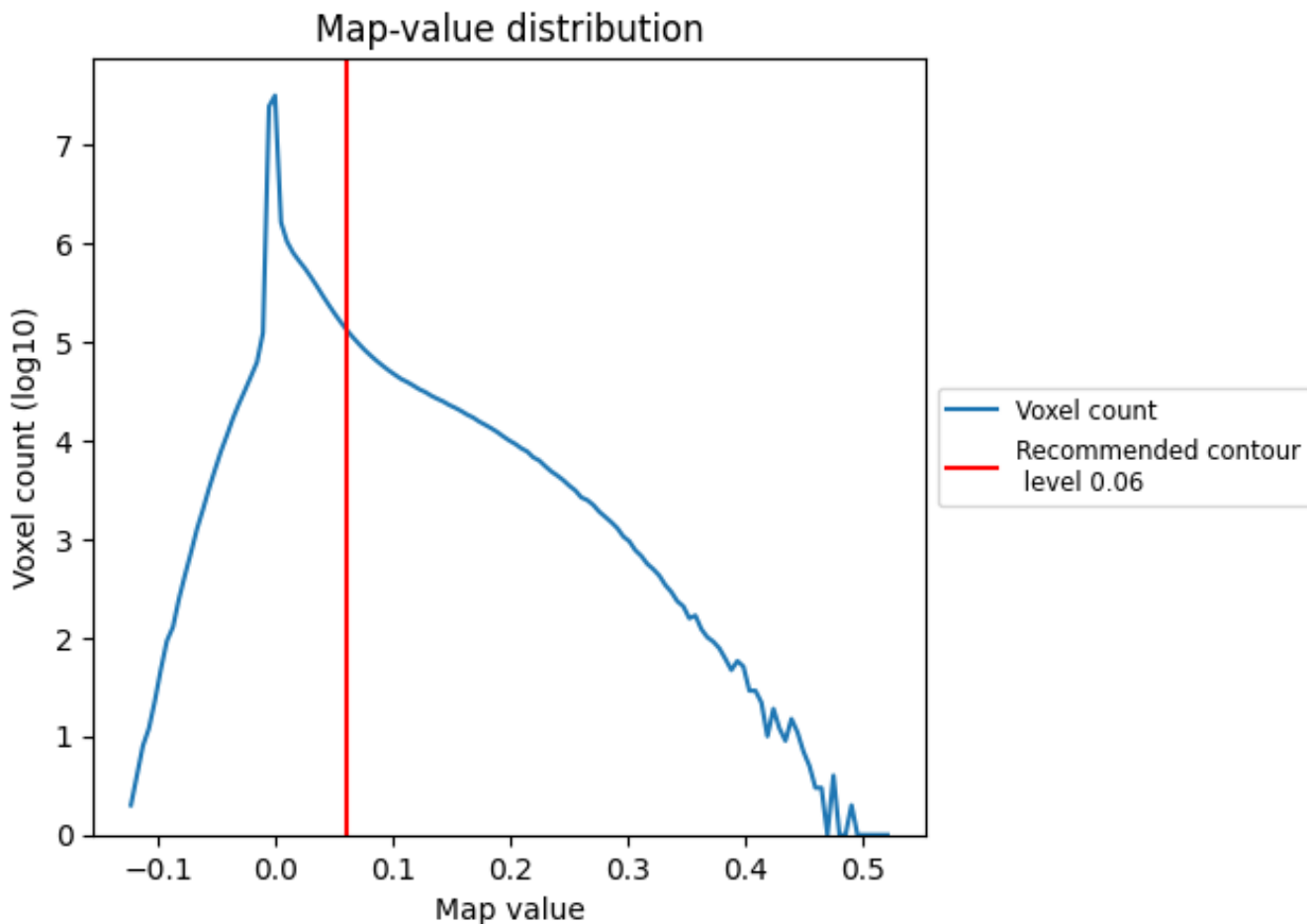
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

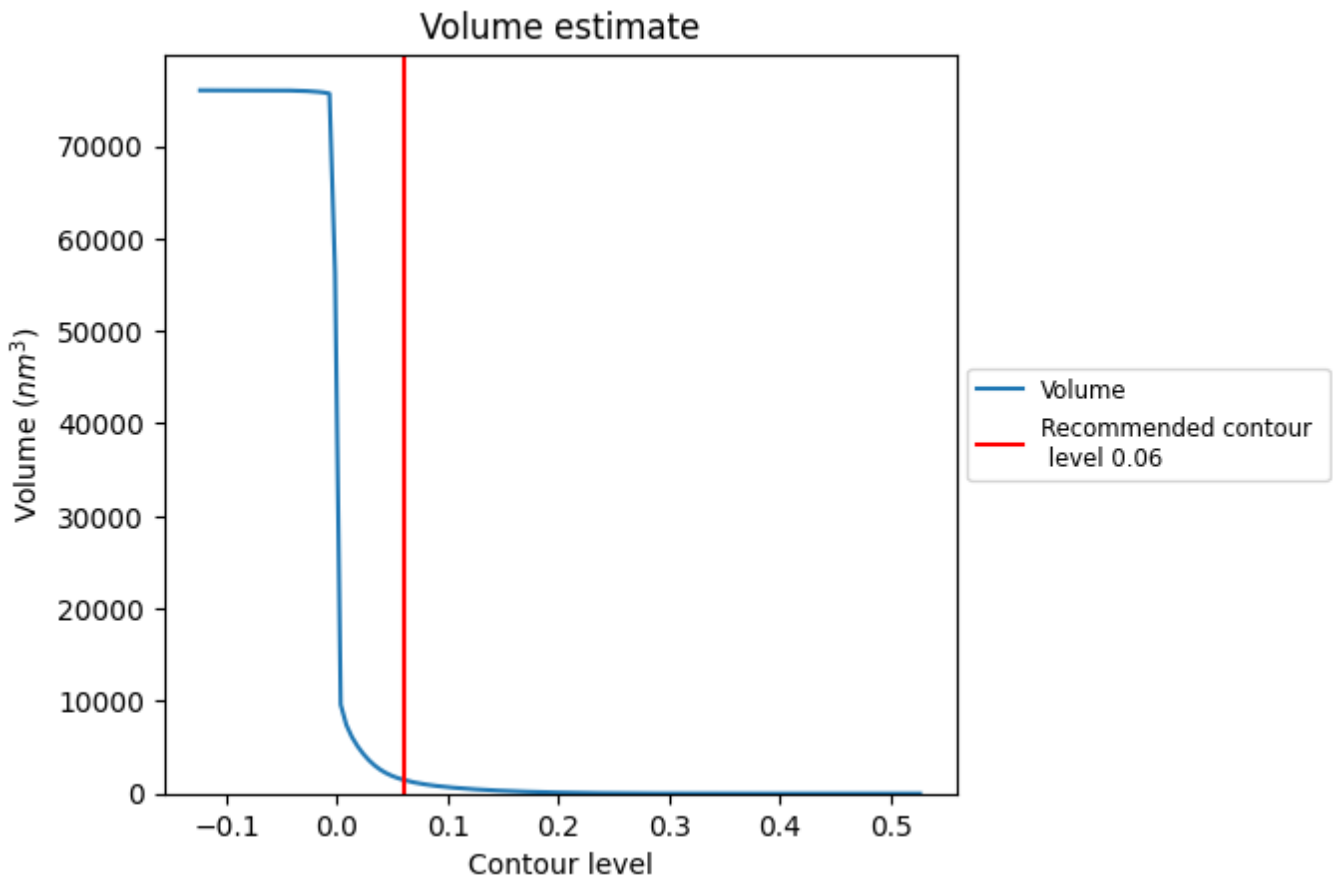
This section contains the results of statistical analysis of the map.

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



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

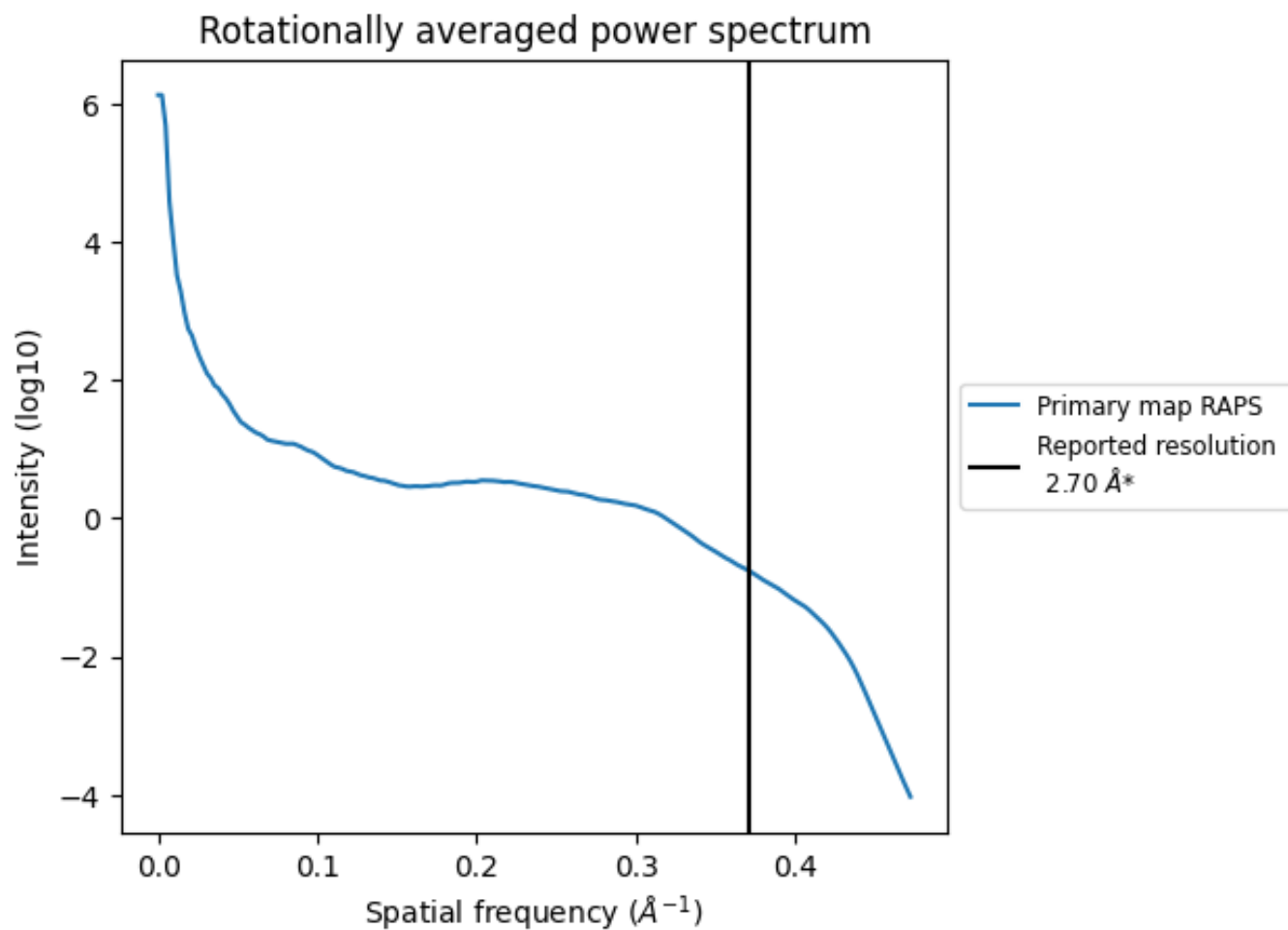
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1485 nm<sup>3</sup>; this corresponds to an approximate mass of 1342 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

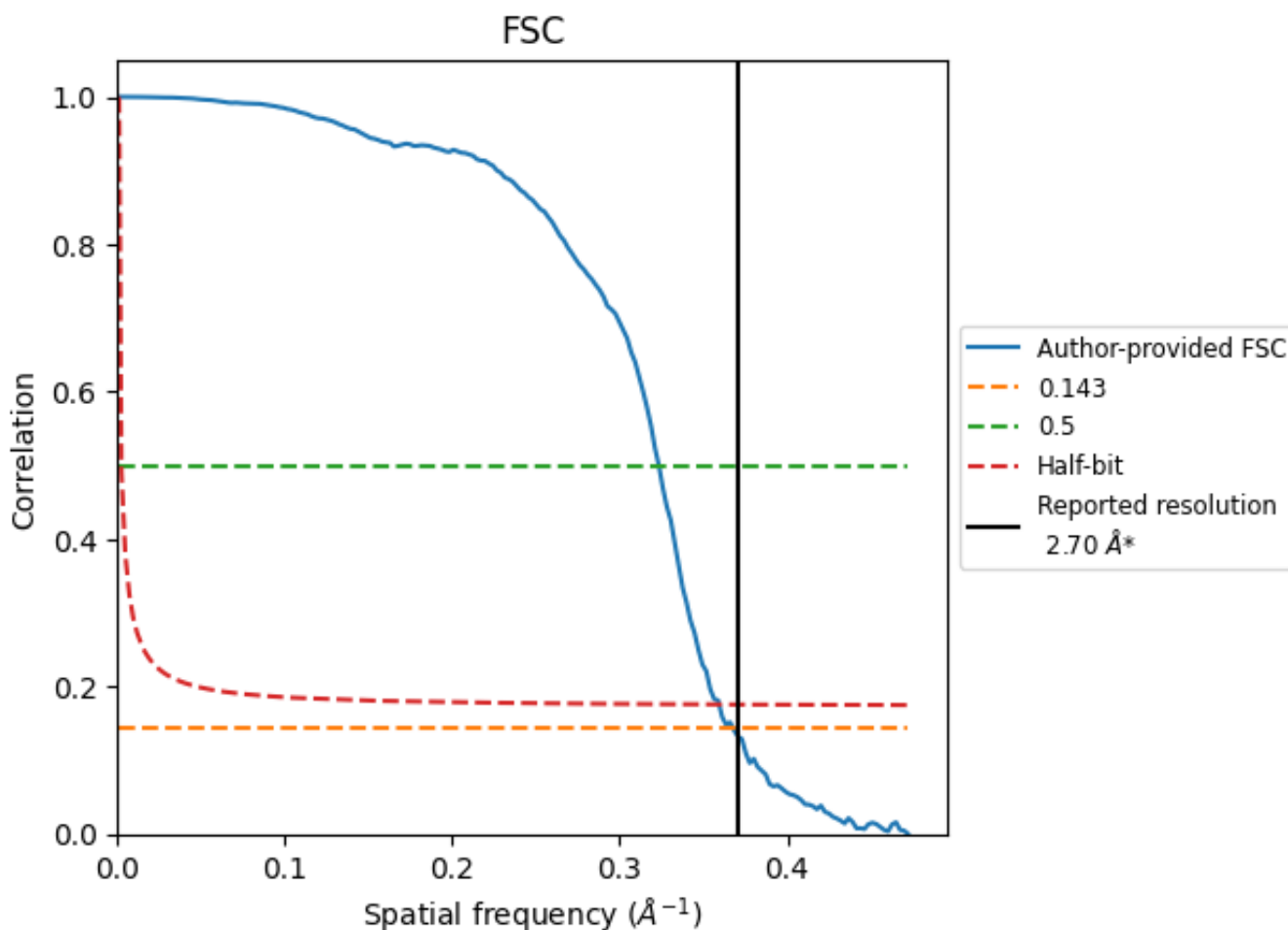


\*Reported resolution corresponds to spatial frequency of 0.370 Å<sup>-1</sup>

## 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 0.370 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

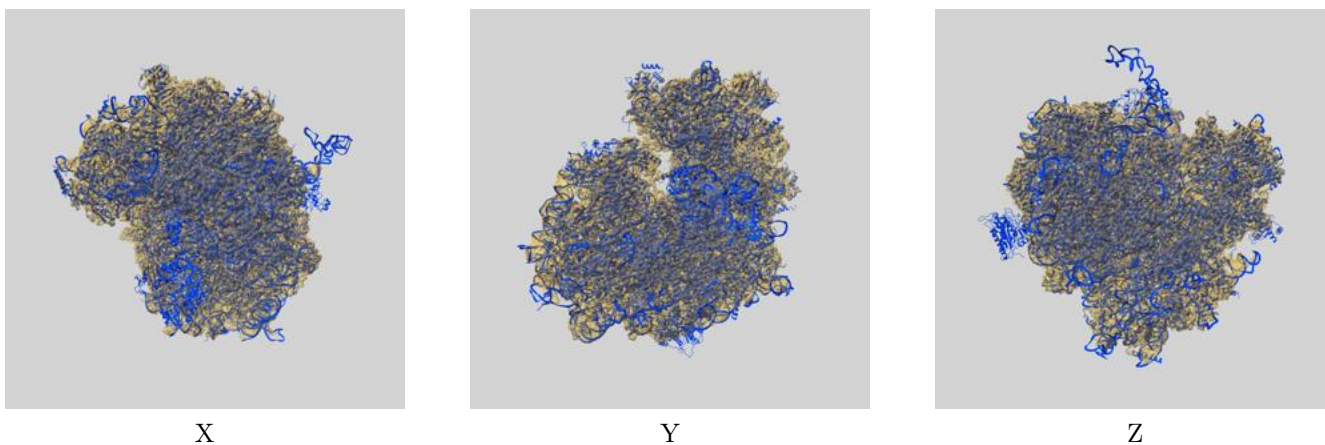
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	2.72	3.09	2.78
Unmasked-calculated*	-	-	-

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

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

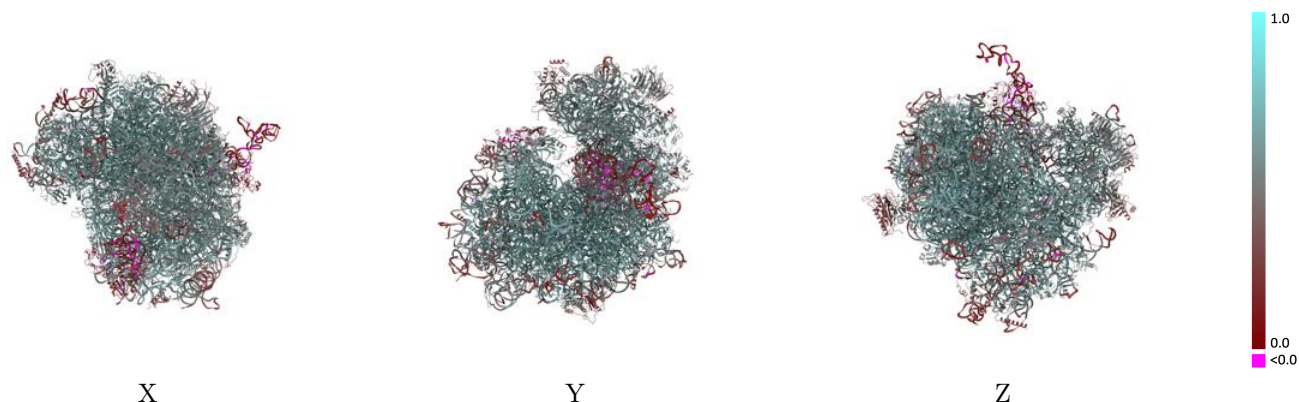
This section contains information regarding the fit between EMDB map EMD-11288 and PDB model 6ZM7. Per-residue inclusion information can be found in section 3 on page 21.

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



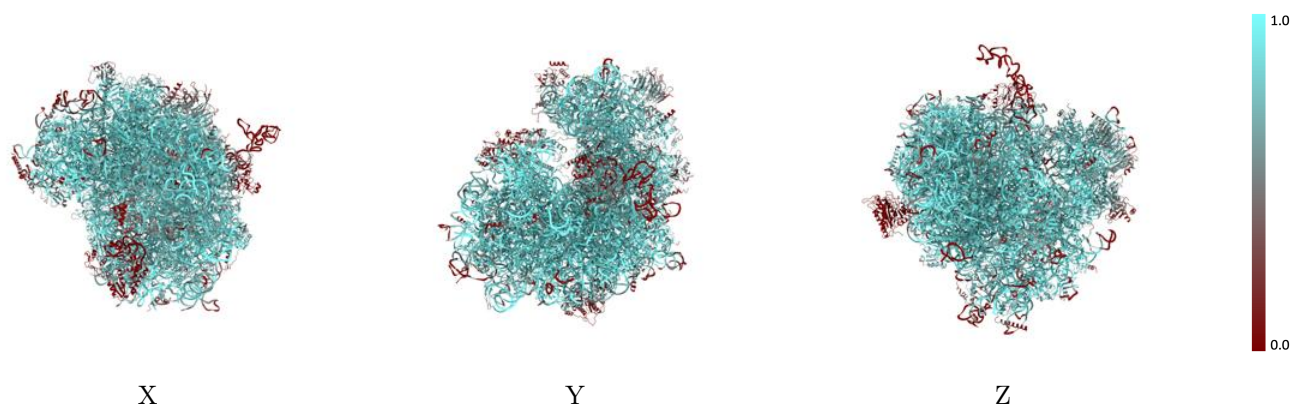
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

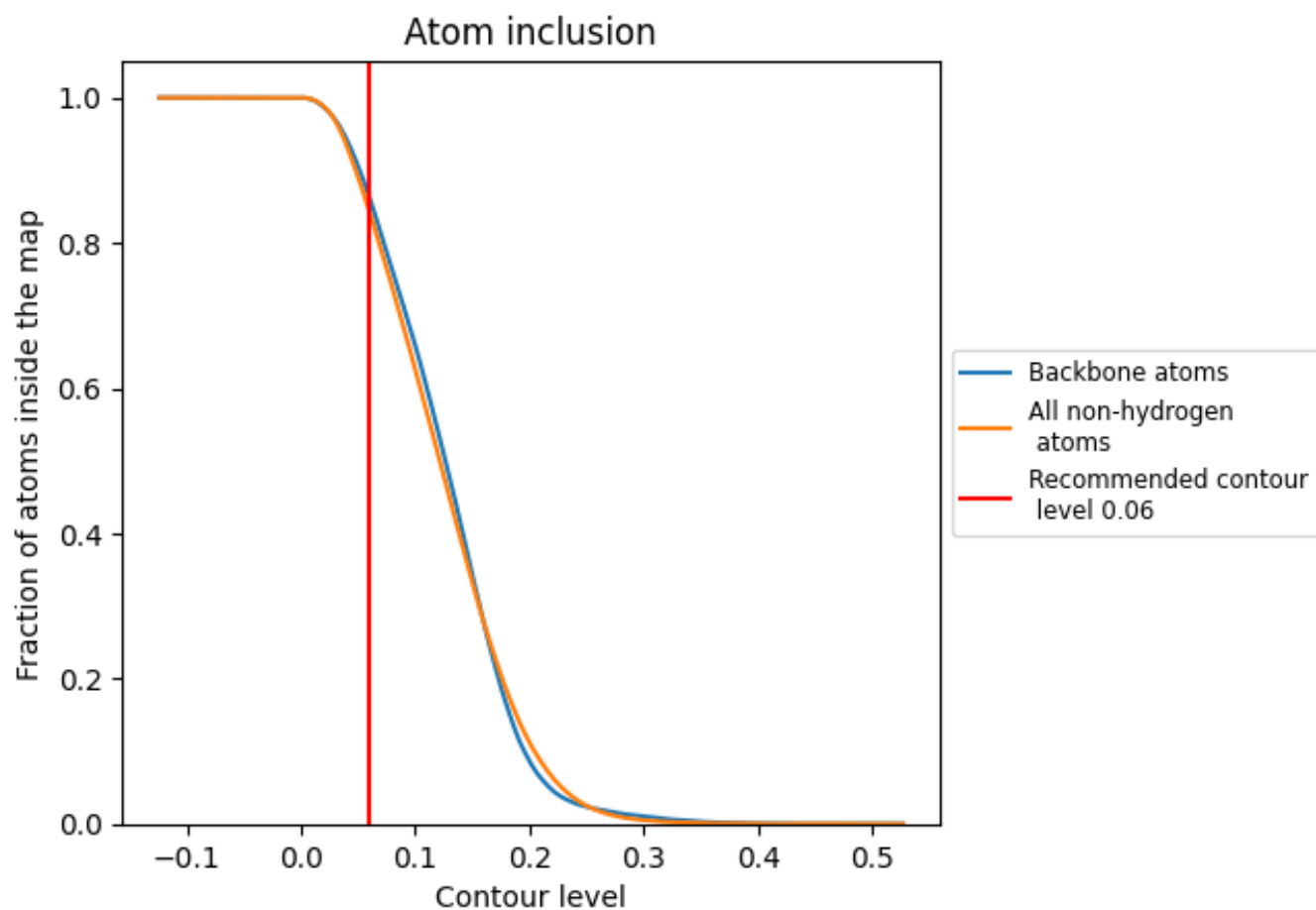
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.06).









































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 84% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8433	 0.5630
CA	 0.0183	 0.3060
CC	 0.7319	 0.3280
CE	 0.5486	 0.4480
CF	 0.8841	 0.5990
L5	 0.8931	 0.5770
L7	 0.9844	 0.6210
L8	 0.9458	 0.6100
LA	 0.9771	 0.6640
LB	 0.9099	 0.6270
LC	 0.9102	 0.6150
LD	 0.8482	 0.5700
LE	 0.7876	 0.5500
LF	 0.9413	 0.6320
LG	 0.7816	 0.5550
LH	 0.8797	 0.5930
LI	 0.9162	 0.6110
LJ	 0.7378	 0.4940
LL	 0.8611	 0.5910
LM	 0.8853	 0.5830
LN	 0.9809	 0.6660
LO	 0.9342	 0.6370
LP	 0.9412	 0.6450
LQ	 0.9620	 0.6500
LR	 0.8512	 0.5920
LS	 0.9507	 0.6380
LT	 0.8979	 0.6050
LU	 0.7423	 0.5140
LV	 0.9426	 0.6360
LW	 0.5949	 0.4730
LX	 0.9050	 0.6150
LY	 0.8729	 0.6060
LZ	 0.8779	 0.6050
La	 0.9487	 0.6530
Lb	 0.7898	 0.5570























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Chain	Atom inclusion	Q-score
Lc	 0.8889	 0.5970
Ld	 0.8775	 0.6070
Le	 0.9538	 0.6470
Lf	 0.9632	 0.6460
Lg	 0.9093	 0.6250
Lh	 0.8807	 0.5980
Li	 0.8797	 0.5980
Lj	 0.9762	 0.6610
Lk	 0.7594	 0.5520
Ll	 0.9598	 0.6310
Lm	 0.9062	 0.6050
Ln	 0.9474	 0.6440
Lo	 0.8925	 0.6210
Lp	 0.9289	 0.6480
Lr	 0.9346	 0.6120
Ls	 0.0312	 0.1650
Lt	 0.0019	 0.1070
Lz	 0.0589	 0.1400
S2	 0.9063	 0.5720
SA	 0.8265	 0.5640
SB	 0.8530	 0.5840
SC	 0.8900	 0.5970
SD	 0.7201	 0.5010
SE	 0.8657	 0.5850
SF	 0.8045	 0.5390
SG	 0.6864	 0.4930
SH	 0.6415	 0.4830
SI	 0.8439	 0.5740
SJ	 0.8553	 0.5660
SK	 0.6932	 0.4720
SL	 0.8531	 0.6050
SM	 0.1519	 0.2650
SN	 0.9070	 0.6100
SO	 0.8608	 0.5800
SP	 0.6582	 0.4700
SQ	 0.7866	 0.5490
SR	 0.7170	 0.5140
SS	 0.7476	 0.5110
ST	 0.7726	 0.5310
SU	 0.6663	 0.4830
SV	 0.8376	 0.5700
SW	 0.9406	 0.6220

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Chain	Atom inclusion	Q-score
SX	 0.9141	 0.6140
SY	 0.7536	 0.5160
SZ	 0.6312	 0.4800
Sa	 0.8885	 0.5970
Sb	 0.7856	 0.5700
Sc	 0.7099	 0.5110
Sd	 0.9072	 0.5800
Se	 0.7594	 0.5220
Sf	 0.2715	 0.3270
Sg	 0.5525	 0.4590