



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 04:20 AM UTC

PDB ID : 9NLE / pdb_00009nle
EMDB ID : EMD-29398
Title : E. coli initiation complex with EQ2-EttA in Hydrolytic 1 conformation
Authors : Singh, S.; Hunt, J.F.
Deposited on : 2025-03-03
Resolution : 3.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

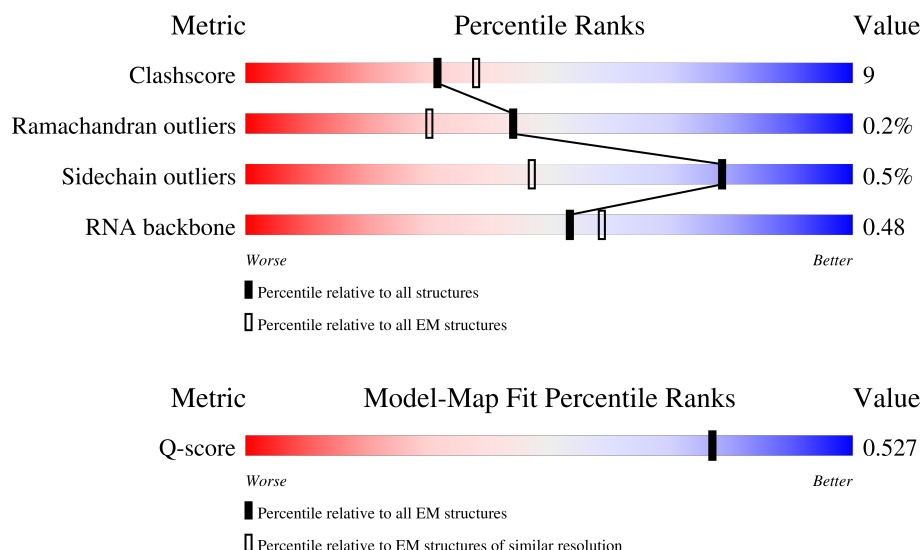
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.20 Å.

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



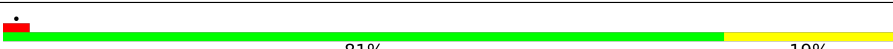
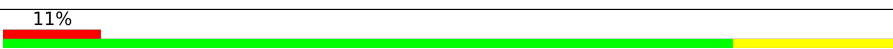
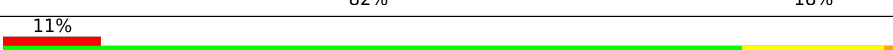
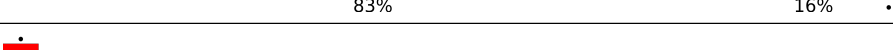
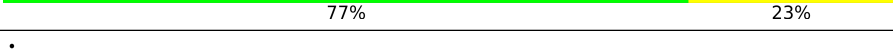
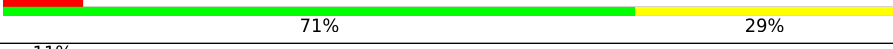
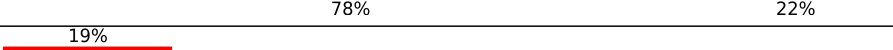
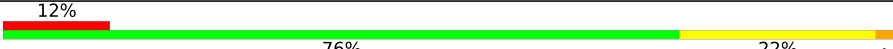
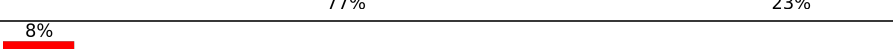
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	15020 (2.70 - 3.70)

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

Mol	Chain	Length	Quality of chain
1	1	220	89% 76% 24%
2	13	142	6% 77% 23%
3	14	122	8% 74% 26%

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Mol	Chain	Length	Quality of chain
4	15	143	
5	16	136	
6	17	120	
7	18	116	
8	19	114	
9	2	271	
10	20	117	
11	21	103	
12	22	110	
13	23	93	
14	24	102	
15	25	94	
16	27	76	
17	28	77	
18	29	63	
19	3	209	
20	30	58	
21	31	66	
22	32	56	
23	33	50	
24	34	46	
25	35	64	
26	36	38	
27	4	201	
28	5	177	

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Mol	Chain	Length	Quality of chain
29	6	176	
30	9	149	
31	E	552	
32	M	9	
33	R1	2903	
34	R2	119	
35	R3	1531	
36	T	77	
37	sb	218	
38	sc	206	
39	sd	205	
40	se	157	
41	sf	100	
42	sg	151	
43	sh	129	
44	si	127	
45	sj	98	
46	sk	116	
47	sl	123	
48	sm	114	
49	sn	100	
50	so	88	
51	sp	82	
52	sq	80	
53	sr	65	

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Mol	Chain	Length	Quality of chain
54	ss	79	
55	st	85	
56	su	70	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
36	H2U	T	20	X	-	-	-
36	4OC	T	32	X	-	-	-
36	MUM	T	54	X	-	-	-
36	PSU	T	55	X	-	-	-
36	4SU	T	8	X	-	-	-

2 Entry composition [i](#)

There are 60 unique types of molecules in this entry. The entry contains 150106 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Large ribosomal subunit protein uL1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	220	Total	C	N	O	S	0	0
			1353	804	270	277	2		

- Molecule 2 is a protein called Large ribosomal subunit protein uL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	13	142	Total	C	N	O	S	0	0
			1129	714	212	199	4		

- Molecule 3 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	14	122	Total	C	N	O	S	0	0
			938	587	180	165	6		

- Molecule 4 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	15	143	Total	C	N	O	S	0	0
			1045	649	206	189	1		

- Molecule 5 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	16	136	Total	C	N	O	S	0	0
			1074	686	205	177	6		

- Molecule 6 is a protein called Large ribosomal subunit protein bL17.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	17	120	Total	C	N	O	S	0	0
			960	593	196	166	5		

- Molecule 7 is a protein called Large ribosomal subunit protein uL18.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	18	116	Total	C	N	O	0	0
			892	552	178	162		

- Molecule 8 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	19	114	Total	C	N	O	S	0	0
			917	574	179	163	1		

- Molecule 9 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	2	271	Total	C	N	O	S	0	0
			2082	1288	423	364	7		

- Molecule 10 is a protein called Large ribosomal subunit protein bL20.

Mol	Chain	Residues	Atoms				AltConf	Trace
10	20	117	Total	C	N	O	0	0
			947	604	192	151		

- Molecule 11 is a protein called Large ribosomal subunit protein bL21.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	21	103	Total	C	N	O	S	0	0
			816	516	153	145	2		

- Molecule 12 is a protein called Large ribosomal subunit protein uL22.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	22	110	Total	C	N	O	S	0	0
			857	532	166	156	3		

- Molecule 13 is a protein called Large ribosomal subunit protein uL23.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	23	93	Total	C	N	O	S	0	0
			738	466	139	131	2		

- Molecule 14 is a protein called Large ribosomal subunit protein uL24.

Mol	Chain	Residues	Atoms				AltConf	Trace
14	24	102	Total	C	N	O		
			779	492	146	141	0	0

- Molecule 15 is a protein called Large ribosomal subunit protein bL25.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	25	94	Total	C	N	O	S		
			753	479	137	134	3	0	0

- Molecule 16 is a protein called Large ribosomal subunit protein bL27.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	27	76	Total	C	N	O	S		
			582	360	117	104	1	0	0

- Molecule 17 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	28	77	Total	C	N	O	S		
			625	388	129	106	2	0	0

- Molecule 18 is a protein called Large ribosomal subunit protein uL29.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	29	63	Total	C	N	O	S		
			509	313	99	95	2	0	0

- Molecule 19 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	3	209	Total	C	N	O	S		
			1565	979	288	294	4	0	0

- Molecule 20 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	30	58	Total	C	N	O	S		
			449	281	87	79	2	0	0

- Molecule 21 is a protein called Large ribosomal subunit protein bL31.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	31	66	Total	C	N	O	S	0	0
			522	323	99	94	6		

- Molecule 22 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	32	56	Total	C	N	O	S	0	0
			444	269	94	80	1		

- Molecule 23 is a protein called Large ribosomal subunit protein bL33.

Mol	Chain	Residues	Atoms				AltConf	Trace
23	33	50	Total	C	N	O	0	0
			409	263	75	71		

- Molecule 24 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	34	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 25 is a protein called Large ribosomal subunit protein bL35.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	35	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 26 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	36	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 27 is a protein called Large ribosomal subunit protein uL4.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	4	201	Total	C	N	O	S	0	0
			1552	974	283	290	5		

- Molecule 28 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	5	177	Total	C	N	O	S	0	0
			1410	899	249	256	6		

- Molecule 29 is a protein called Large ribosomal subunit protein uL6.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	6	176	Total	C	N	O	S	0	0
			1323	832	243	246	2		

- Molecule 30 is a protein called Large ribosomal subunit protein bL9.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	9	149	Total	C	N	O	S	0	0
			1111	699	197	214	1		

- Molecule 31 is a protein called Energy-dependent translational throttle protein EttA.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	E	552	Total	C	N	O	S	0	0
			4360	2745	774	829	12		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	188	GLN	GLU	conflict	UNP P0A9W3
E	341	SER	ASP	conflict	UNP P0A9W3
E	470	GLN	GLU	conflict	UNP P0A9W3

- Molecule 32 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	M	9	Total	C	N	O	P	0	0
			195	88	40	58	9		

- Molecule 33 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	R1	2903	Total	C	N	O	P	0	0
			62318	27801	11467	20148	2902		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R1	1847	G	A	conflict	GB 2019144442

- Molecule 34 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	R2	119	Total	C	N	O	P	0	0
			2546	1135	466	827	118		

- Molecule 35 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	R3	1531	Total	C	N	O	P	0	0
			32850	14652	6028	10640	1530		

- Molecule 36 is a RNA chain called tRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
36	T	77	Total	C	N	O	P	S	0	0
			1639	734	294	534	76	1		

- Molecule 37 is a protein called Small ribosomal subunit protein uS2.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	sb	218	Total	C	N	O	S	0	0
			1704	1081	305	311	7		

- Molecule 38 is a protein called Small ribosomal subunit protein uS3.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	sc	206	Total	C	N	O	S	0	0
			1624	1028	305	288	3		

- Molecule 39 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	sd	205	Total	C	N	O	S	0	0
			1643	1026	315	298	4		

- Molecule 40 is a protein called Small ribosomal subunit protein uS5.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	se	157	Total	C	N	O	S	0	0
			1156	719	218	213	6		

- Molecule 41 is a protein called 30S ribosomal protein S6, non-modified isoform.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	sf	100	Total	C	N	O	S	0	0
			817	515	148	148	6		

- Molecule 42 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	sg	151	Total	C	N	O	S	0	0
			1181	735	227	215	4		

- Molecule 43 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	sh	129	Total	C	N	O	S	0	0
			979	616	173	184	6		

- Molecule 44 is a protein called Small ribosomal subunit protein uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	si	127	Total	C	N	O	S	0	0
			1022	634	206	179	3		

- Molecule 45 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	sj	98	Total	C	N	O	S	0	0
			786	493	150	142	1		

- Molecule 46 is a protein called Small ribosomal subunit protein uS11.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	sk	116	Total	C	N	O	S	0	0
			869	535	173	158	3		

- Molecule 47 is a protein called Small ribosomal subunit protein uS12.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	sl	123	Total	C	N	O	S	0	0
			955	590	196	165	4		

- Molecule 48 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	sm	114	Total	C	N	O	S	0	0
			883	546	178	156	3		

- Molecule 49 is a protein called Small ribosomal subunit protein uS14.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	sn	100	Total	C	N	O	S	0	0
			805	499	164	139	3		

- Molecule 50 is a protein called Small ribosomal subunit protein uS15.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	so	88	Total	C	N	O	S	0	0
			714	439	144	130	1		

- Molecule 51 is a protein called Small ribosomal subunit protein bS16.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	sp	82	Total	C	N	O	S	0	0
			649	406	128	114	1		

- Molecule 52 is a protein called Small ribosomal subunit protein uS17.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	sq	80	Total	C	N	O	S	0	0
			648	411	121	113	3		

- Molecule 53 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	sr	65	Total	C	N	O	S	0	0
			535	339	100	95	1		

- Molecule 54 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	ss	79	Total	C	N	O	S	0	0
			637	408	120	107	2		

- Molecule 55 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	st	85	Total	C	N	O	S	0	0
			665	411	137	114	3		

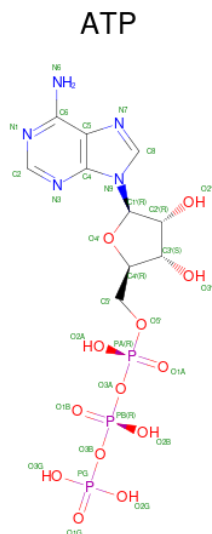
- Molecule 56 is a protein called Small ribosomal subunit protein bS21.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	su	70	Total	C	N	O	S	0	0
			590	366	125	98	1		

- Molecule 57 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
57	15	1	Total	Mg	0
			1	1	
57	17	1	Total	Mg	0
			1	1	
57	M	1	Total	Mg	0
			1	1	
57	R1	204	Total	Mg	0
			204	204	
57	R3	89	Total	Mg	0
			89	89	
57	sn	1	Total	Mg	0
			1	1	
57	sq	1	Total	Mg	0
			1	1	

- Molecule 58 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).

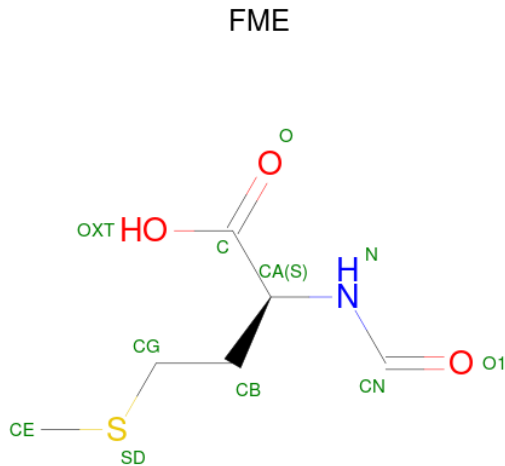


Mol	Chain	Residues	Atoms					AltConf
58	E	1	Total 31	C 10	N 5	O 13	P 3	0
58	E	1	Total 31	C 10	N 5	O 13	P 3	0

- Molecule 59 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		AltConf
59	E	2	Total 2	Na 2	0

- Molecule 60 is N-FORMYLMETHIONINE (CCD ID: FME) (formula: $\text{C}_6\text{H}_{11}\text{NO}_3\text{S}$).

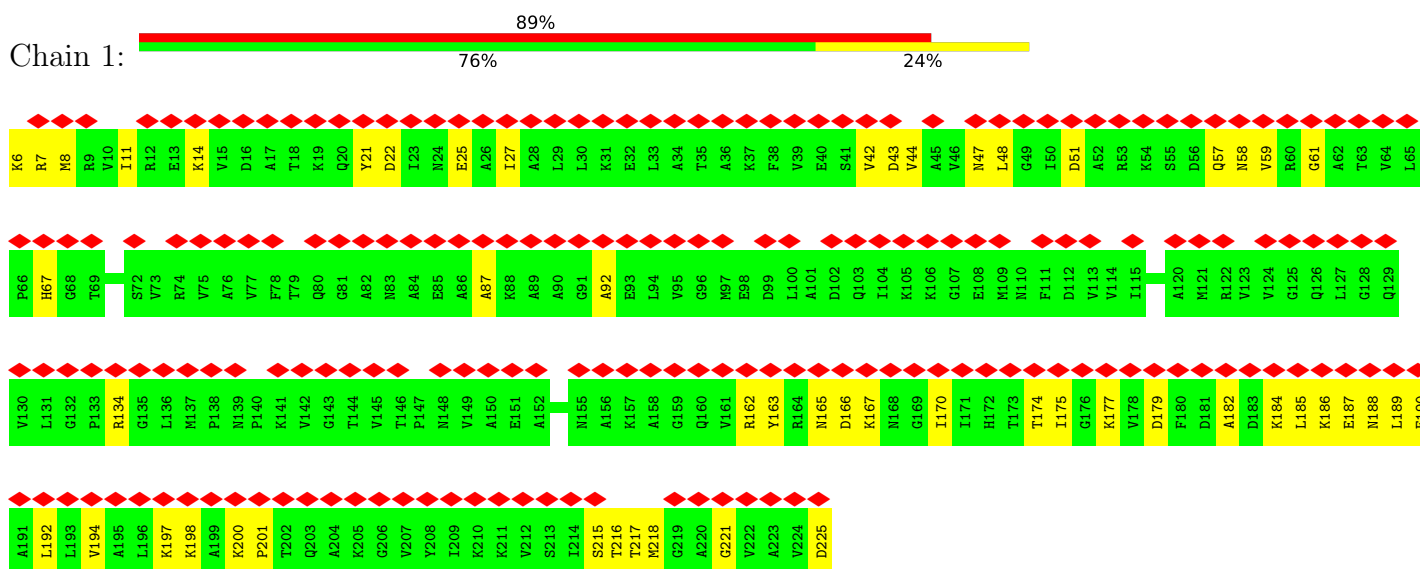


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	S	
60	T	1	10	6	1	2	1	0

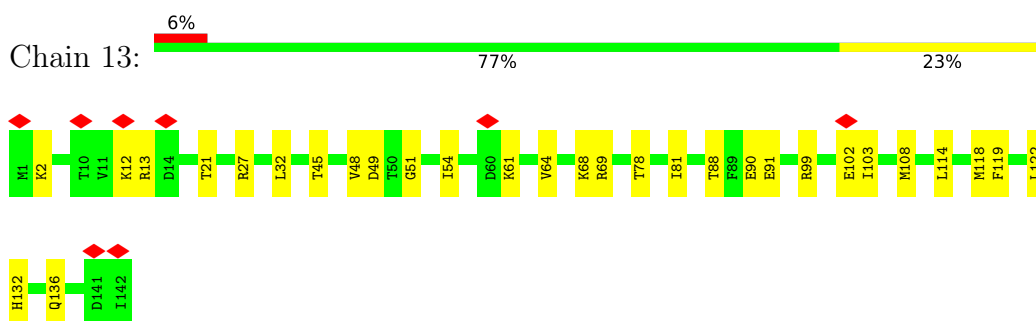
3 Residue-property plots

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

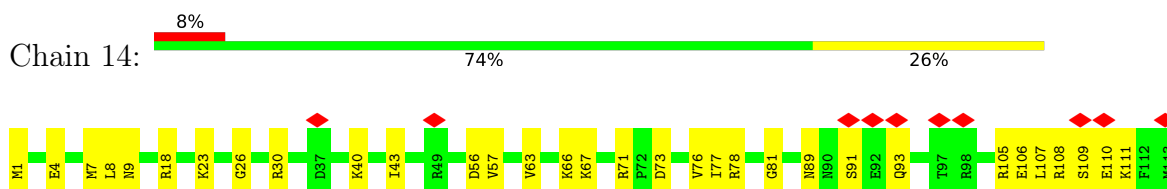
- Molecule 1: Large ribosomal subunit protein uL1



- Molecule 2: Large ribosomal subunit protein uL13



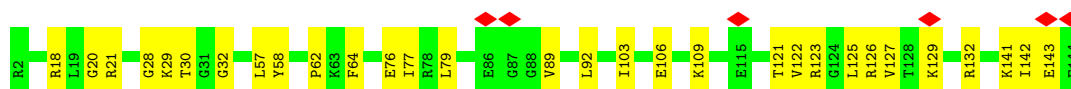
- Molecule 3: 50S ribosomal protein L14





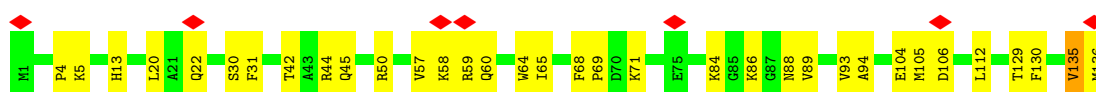
- Molecule 4: 50S ribosomal protein L15

Chain 15: 79% 21%



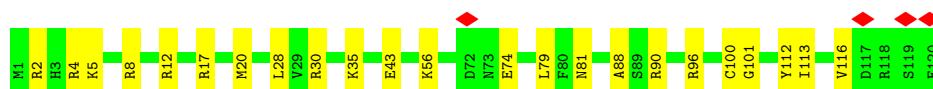
- Molecule 5: 50S ribosomal protein L16

Chain 16: 5% 75% 24%



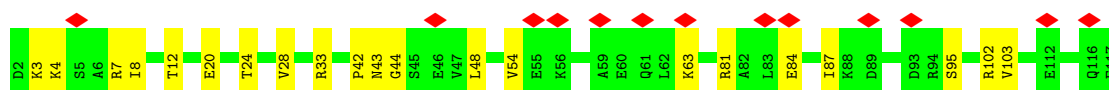
- Molecule 6: Large ribosomal subunit protein bL17

Chain 17: 81% 19%



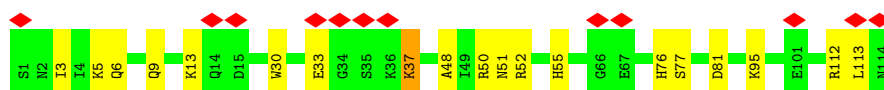
- Molecule 7: Large ribosomal subunit protein uL18

Chain 18: 11% 82% 18%



- Molecule 8: 50S ribosomal protein L19

Chain 19: 11% 83% 16%



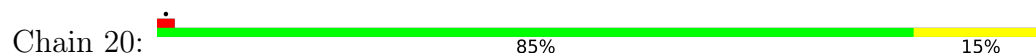
- Molecule 9: 50S ribosomal protein L2

Chain 2: 77% 23%





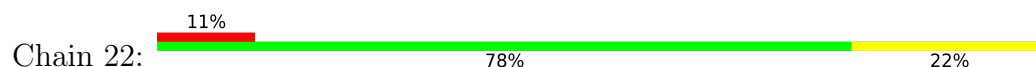
- Molecule 10: Large ribosomal subunit protein bL20



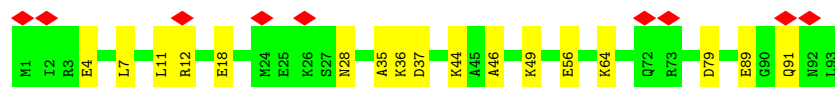
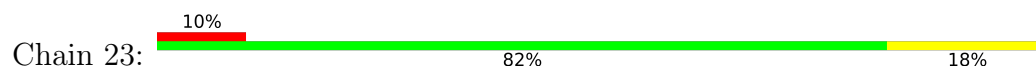
- Molecule 11: Large ribosomal subunit protein bL21



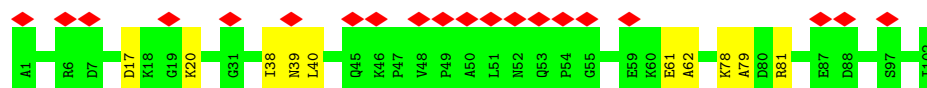
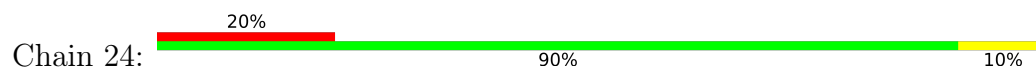
- Molecule 12: Large ribosomal subunit protein uL22



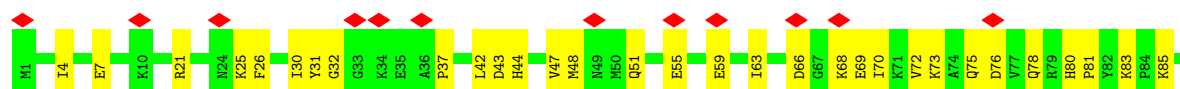
- Molecule 13: Large ribosomal subunit protein uL23

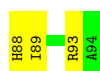


- Molecule 14: Large ribosomal subunit protein uL24

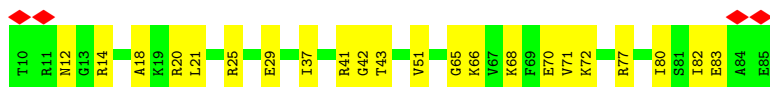
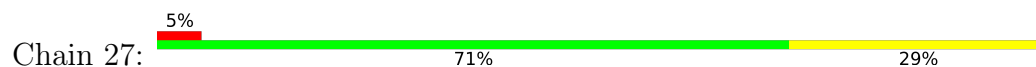


- Molecule 15: Large ribosomal subunit protein bL25

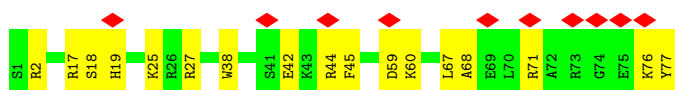
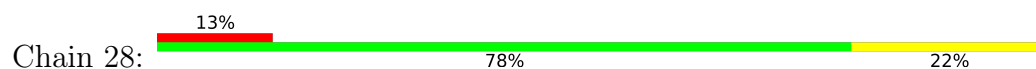




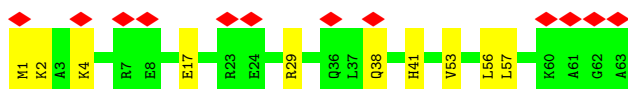
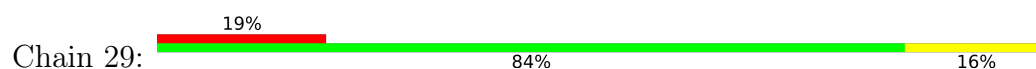
- Molecule 16: Large ribosomal subunit protein bL27



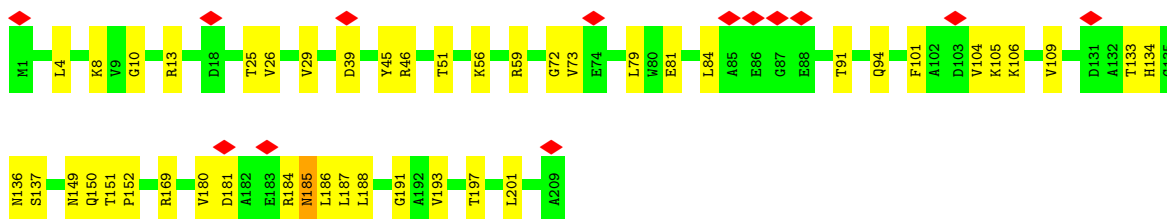
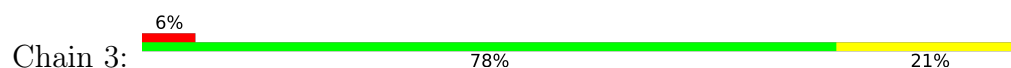
- Molecule 17: 50S ribosomal protein L28



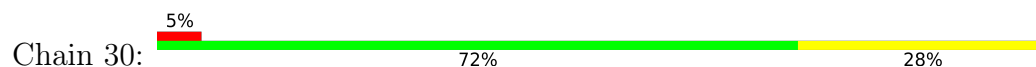
- Molecule 18: Large ribosomal subunit protein uL29



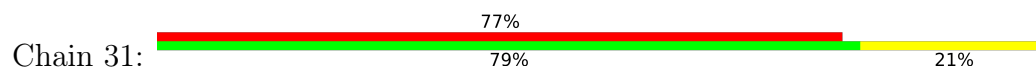
- Molecule 19: 50S ribosomal protein L3

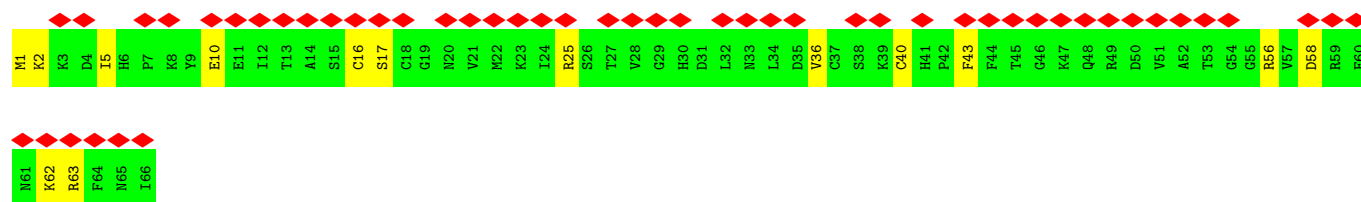


- Molecule 20: 50S ribosomal protein L30

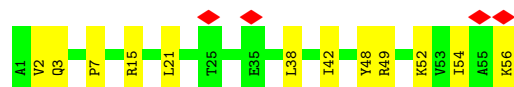
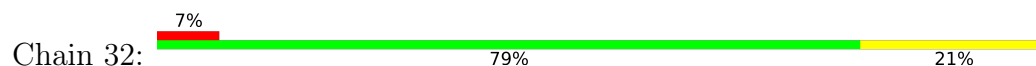


- Molecule 21: Large ribosomal subunit protein bL31

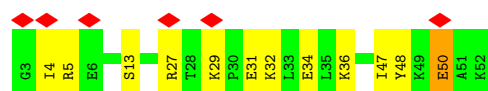
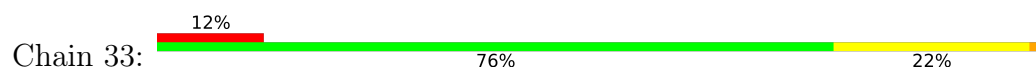




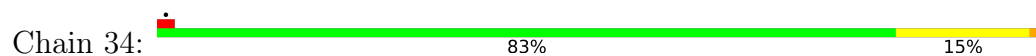
- Molecule 22: 50S ribosomal protein L32



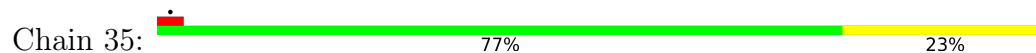
- Molecule 23: Large ribosomal subunit protein bL33



- Molecule 24: 50S ribosomal protein L34



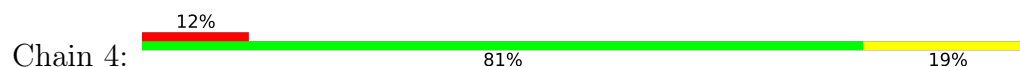
- Molecule 25: Large ribosomal subunit protein bL35

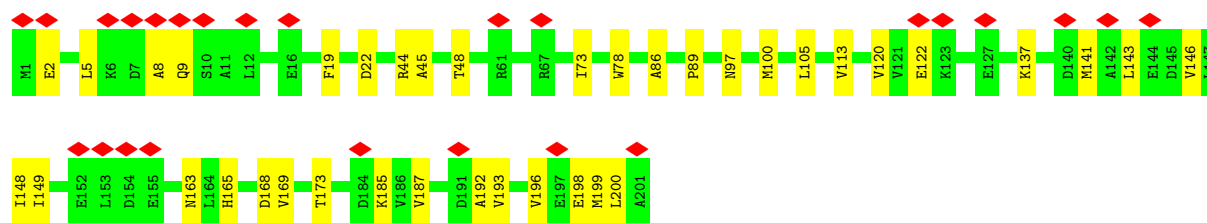


- Molecule 26: 50S ribosomal protein L36

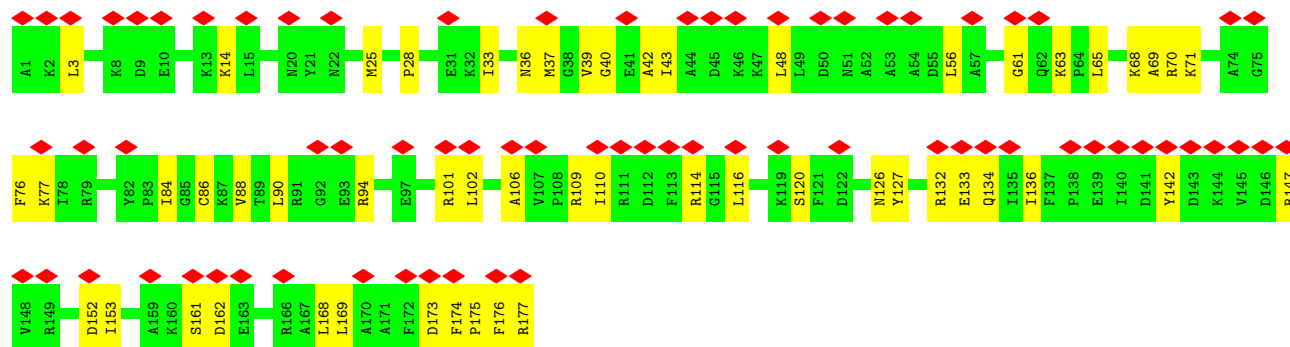
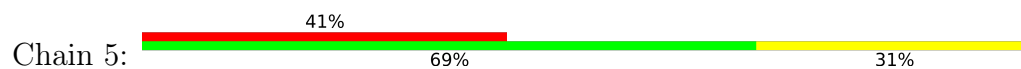


- Molecule 27: Large ribosomal subunit protein uL4

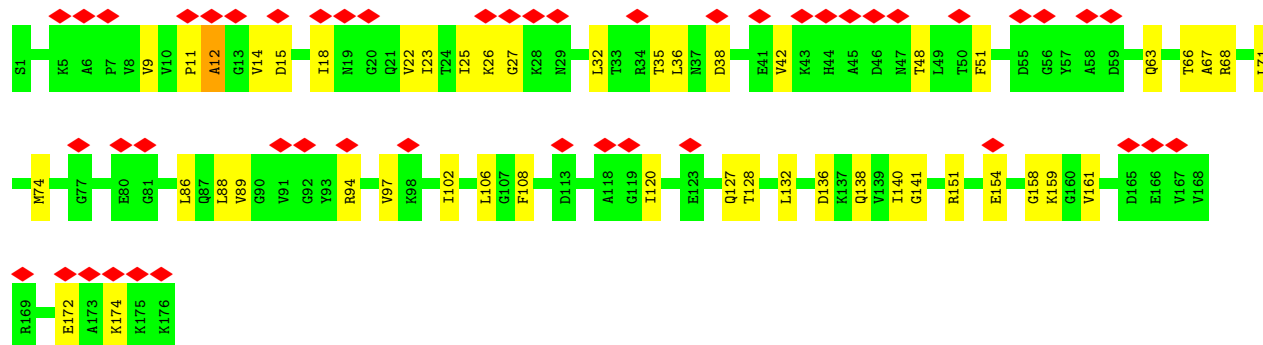
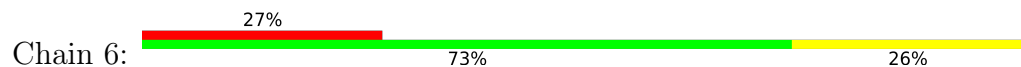




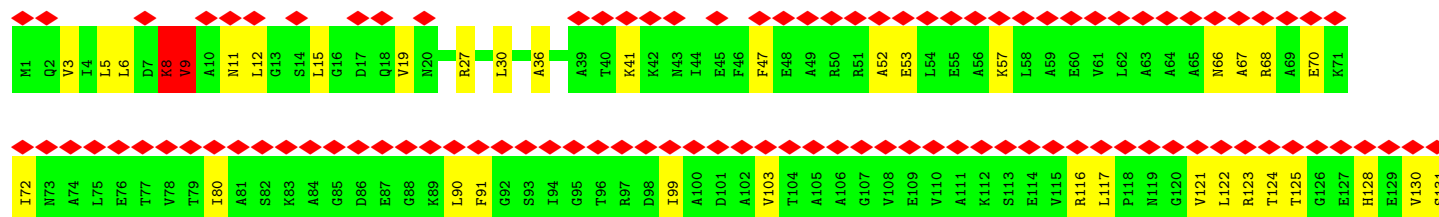
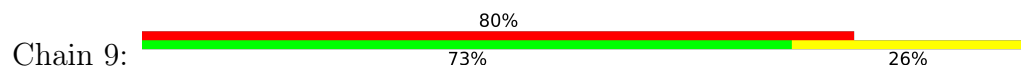
- Molecule 28: 50S ribosomal protein L5

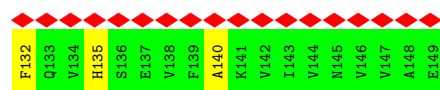


- Molecule 29: Large ribosomal subunit protein uL6

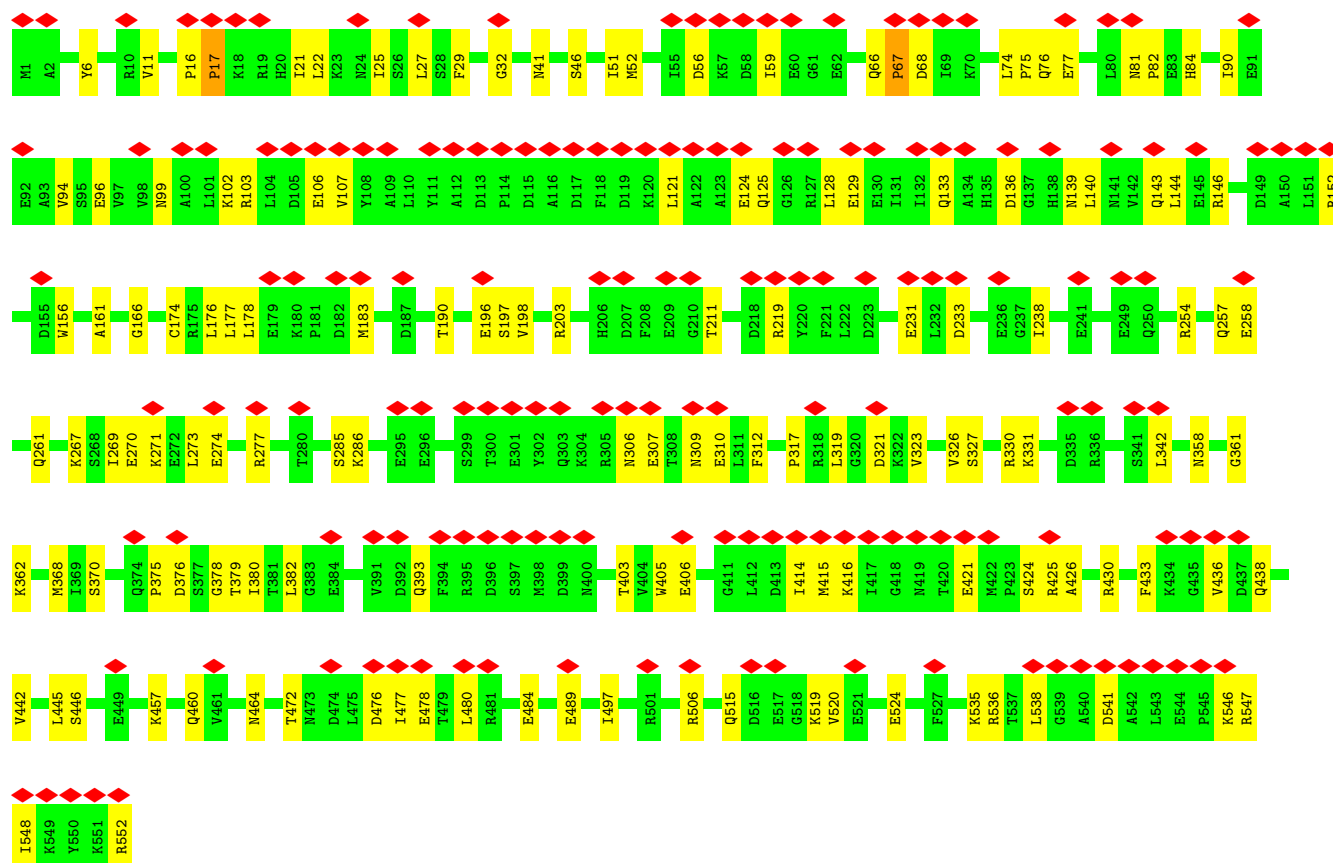
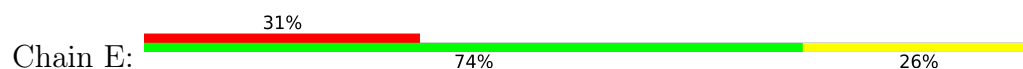


- Molecule 30: Large ribosomal subunit protein bL9





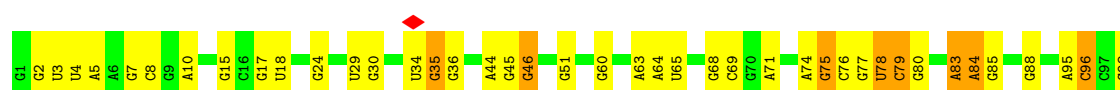
• Molecule 31: Energy-dependent translational throttle protein EttA

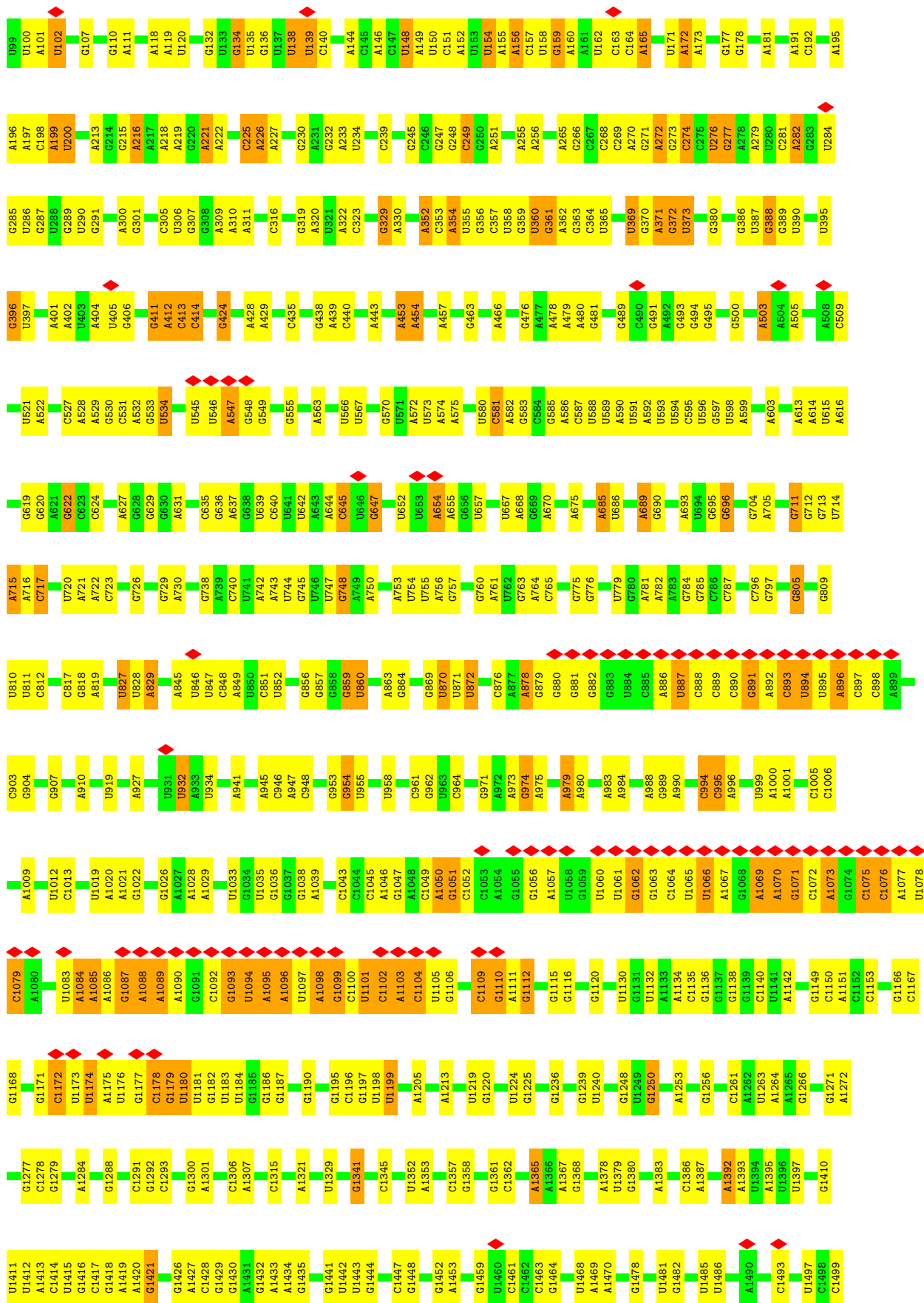


• Molecule 32: mRNA

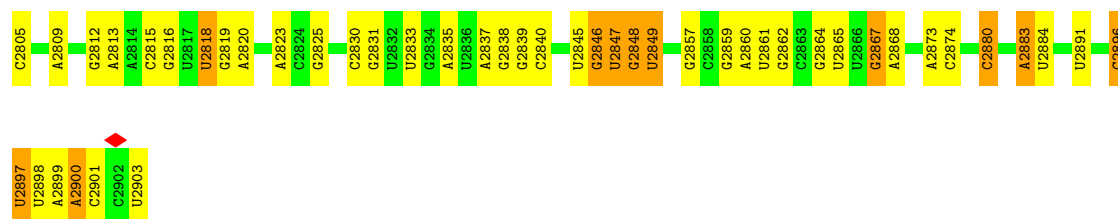


• Molecule 33: 23S ribosomal RNA

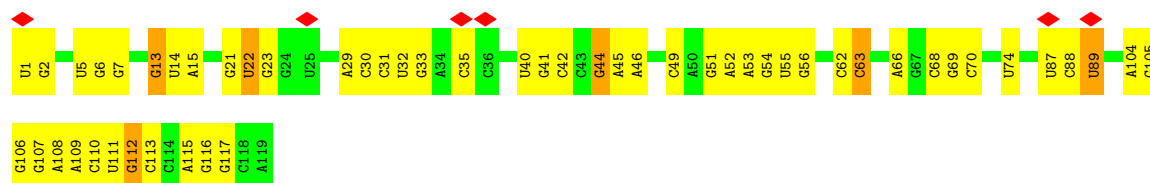




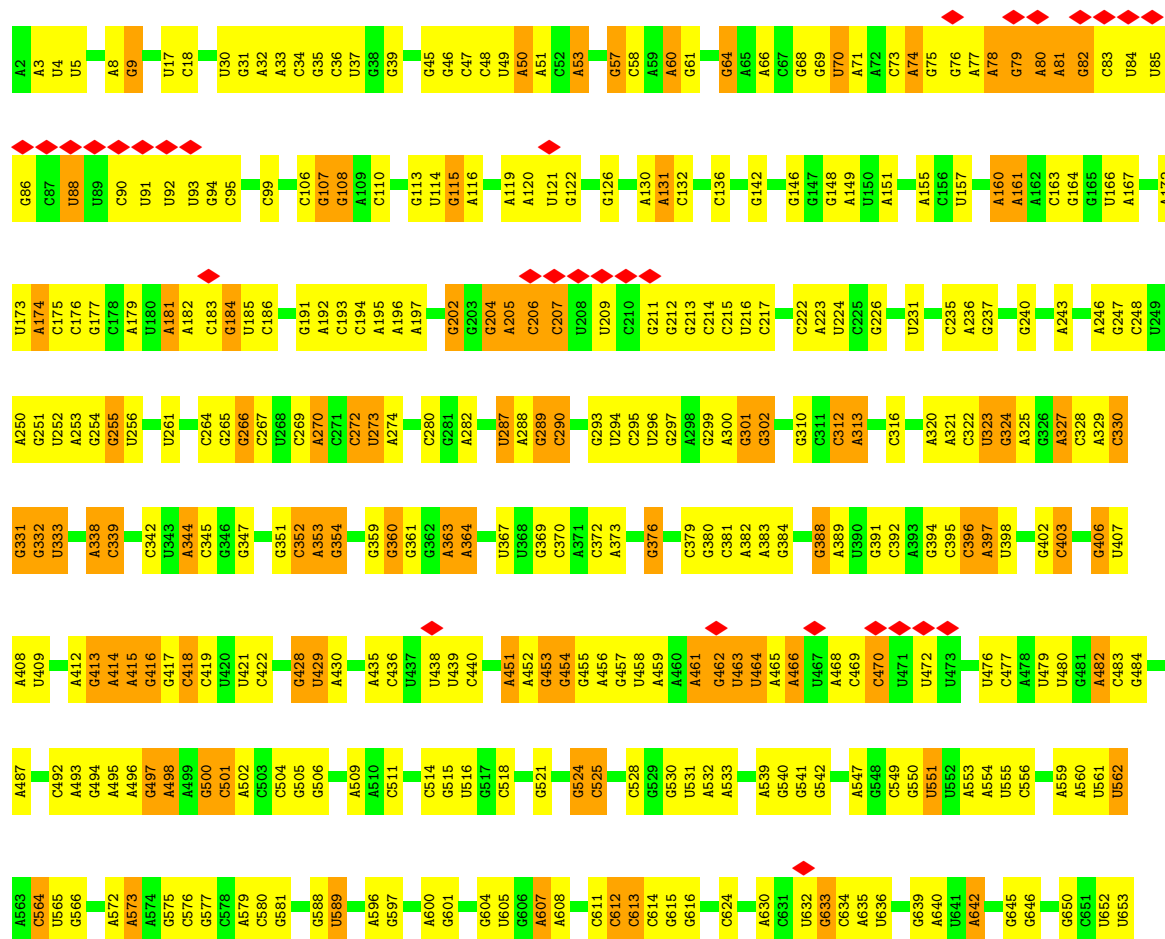
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G2581	G2582	C2591	C2592	A2598	U2477	A2602	U2609	U2613	A2614	U2615	G2623	G2624	G2625	G2626	U2629	U2630	A2635	C2636	U2637	G2638	G2641	G2645	C2646	C2649	U2650	C2651	C2652	U2656	A2662	G2663	G2664	A2665	C2666	C2667	G2668	C2669	A2670	G2671	A2682	G2685	U2689	U2690	G2697								
A2469	G2470	U2474	C2475	A2476	U2477	A2478	G2484	G2494	G2495	C2496	A2497	G2502	A2503	U2504	G2505	U2506	C2507	G2508	C2512	A2513	U2514	C2515	A2516	C2517	U2518	U2519	C2520	G2529	U2537	C2538	G2545	U2546	A2547	U2548	U2554	A2560	U2561	U2562	U2563	A2564	A2565	A2566	U2567	U2568	A2572	C2579	U2580				
G2382	G2383	U2384	C2385	A2386	G2389	U2390	G2391	G2396	G2400	U2401	U2402	A2406	A2407	U2408	G2409	G2410	A2411	A2412	G2415	C2416	C2417	U2423	C2424	A2425	U2426	C2427	G2428	G2429	A2430	A2434	A2435	G2436	A2439	C2440	U2441	C2442	C2443	A2448	G2455	C2456	U2457	A2461	C2462	G2463	G2464	C2467	A2468				
A2284	C2285	G2286	A2287	A2288	G2289	U2290	U2291	U2292	A2297	G2304	U2305	C2306	C2307	G2308	A2309	U2312	C2313	G2314	U2315	G2316	U2321	A2322	C2323	U2324	G2325	C2326	A2327	U2328	U2329	C2330	G2331	C2332	A2333	G2345	U2346	C2347	U2348	G2349	C2350	U2356	G2361	G2365	U2371	U2372	C2373	C2374	C2375	U2376	A2377	C2380	A2381
U2185	G2186	U2187	U2188	U2189	G2190	A2191	U2192	U2193	U2194	A2198	U2203	G2204	C2208	U2213	C2214	G2215	U2216	G2217	U2218	U2219	G2224	A2225	U2226	U2229	G2230	U2233	C2234	G2235	U2236	U2237	G2238	G2239	U2243	U2244	U2245	G2246	C2248	U2249	G2250	C2258	A2266	G2271	U2272	A2273	C2282	C2283					
C2043	C2047	G2048	A2051	C2055	G2056	G2057	A2058	A2059	A2060	G2061	A2062	C2066	G2067	U2068	G2069	A2070	A2071	C2072	C2073	A2080	U2081	A2082	U2086	G2087	G2093	C2096	U2099	G2100	A2101	G2102	C2103	C2104	G2107	A2108	U2109	G2110	U2111	G2112	G2115	U2116	A2117	U2118	A2119	G2120	G2121	U2122	G2123				
A1918	A1919	U1923	C1924	A1927	A1928	A1929	G1930	A1936	A1937	A1938	U1943	A1952	U1955	A1966	C1967	A1970	U1971	G1972	U1991	C1996	C1997	C2000	G2001	G2002	C2008	A2009	G2012	A2013	A2014	A2015	U2016	U2022	C2023	G2024	C2025	U2026	A2030	G2032	A2033	U2034	G2038	U2039									
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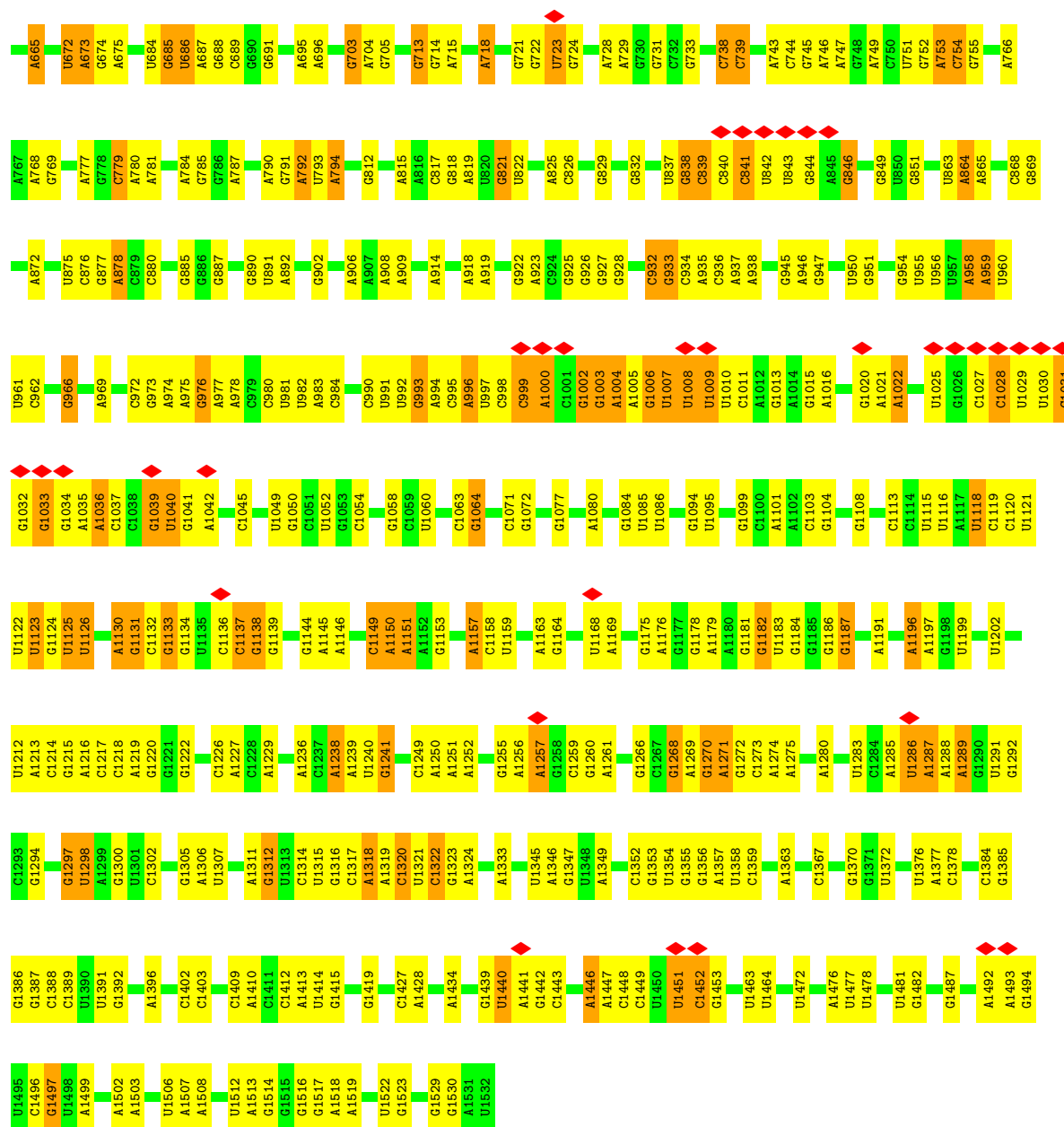


• Molecule 34: 5S ribosomal RNA



• Molecule 35: 16S ribosomal RNA

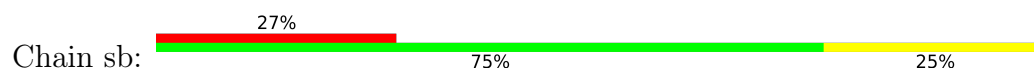


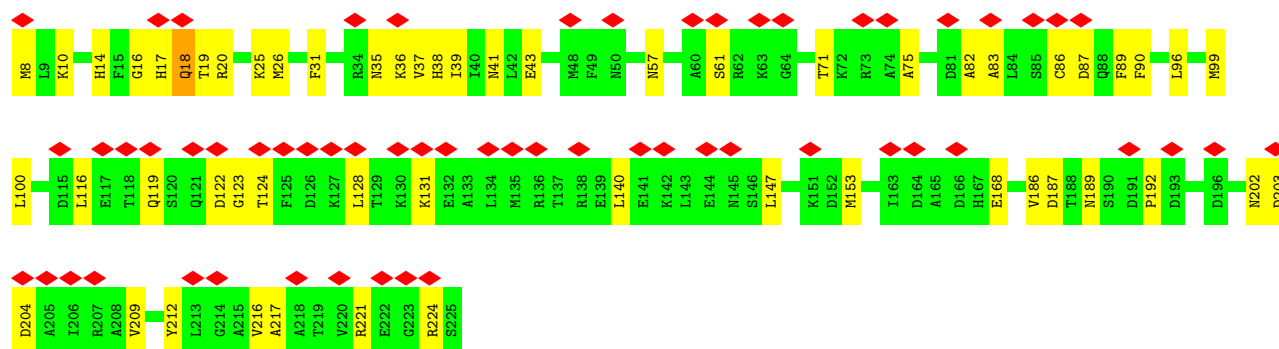


- Molecule 36: tRNA

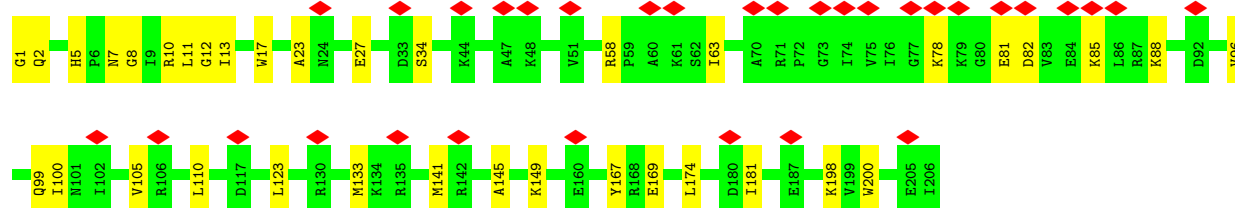
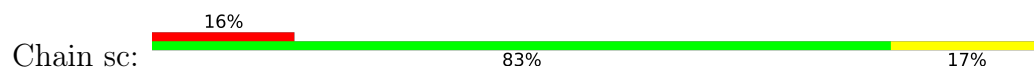


- Molecule 37: Small ribosomal subunit protein uS2

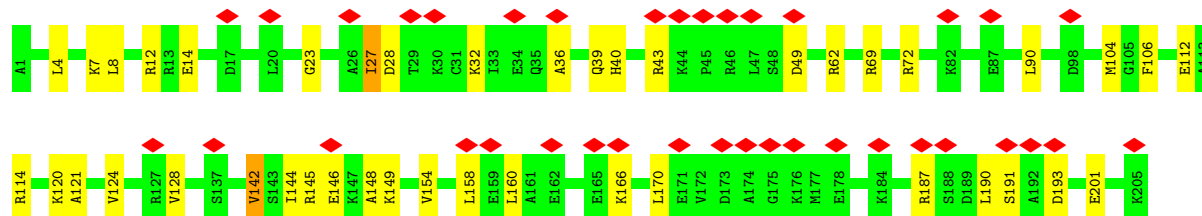
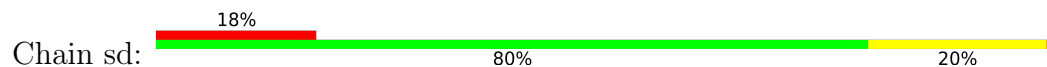




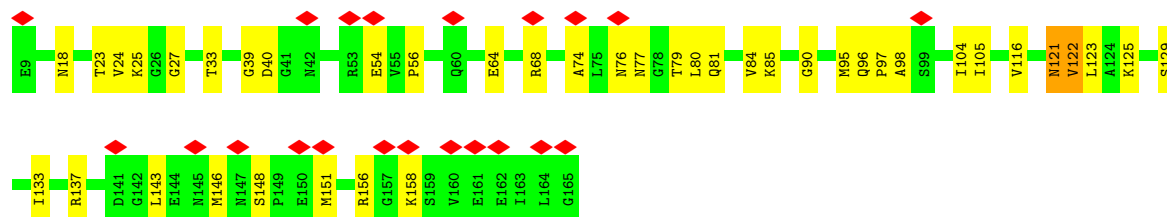
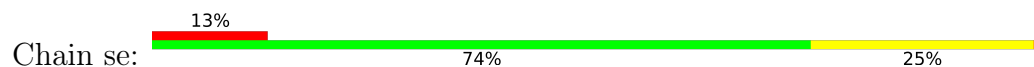
- Molecule 38: Small ribosomal subunit protein uS3



- Molecule 39: 30S ribosomal protein S4

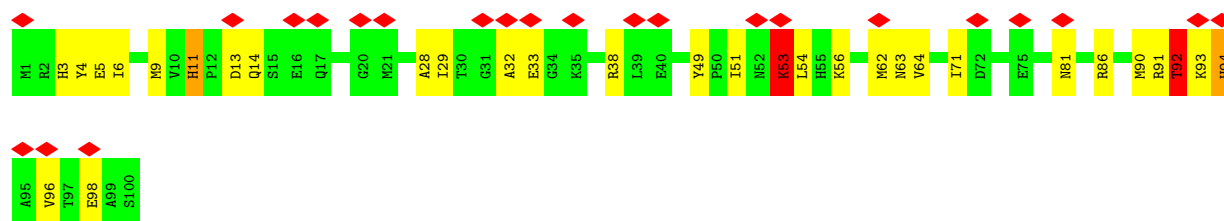


- Molecule 40: Small ribosomal subunit protein uS5

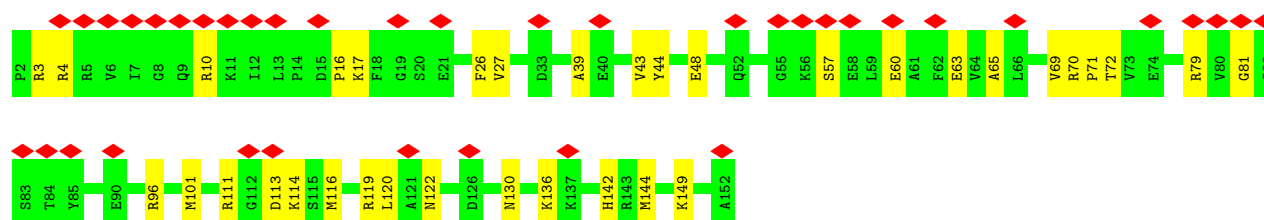
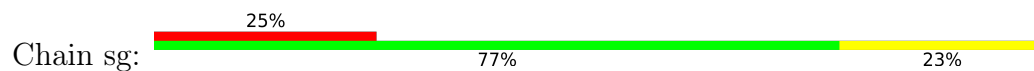


- Molecule 41: 30S ribosomal protein S6, non-modified isoform

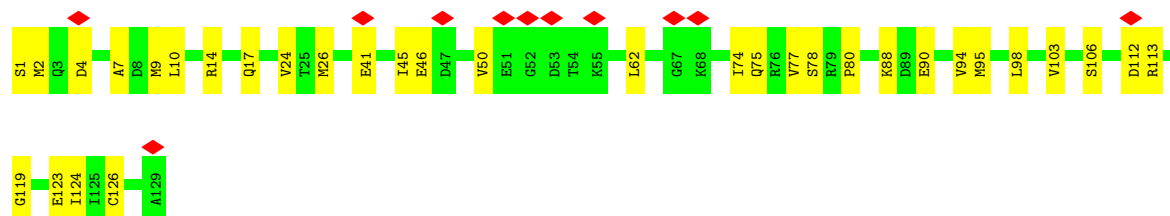
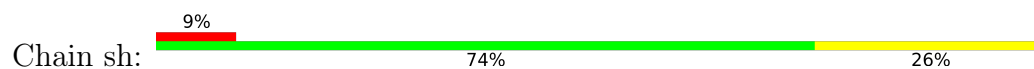




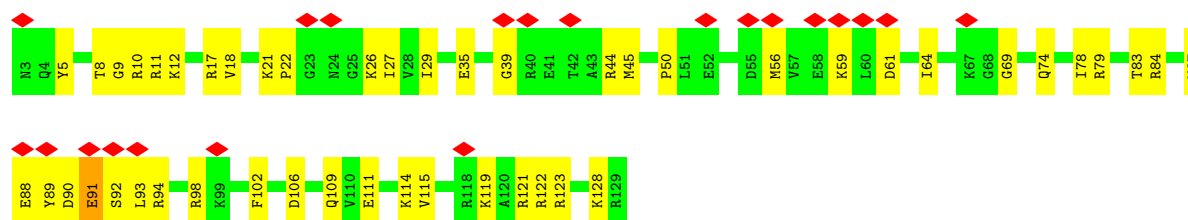
- Molecule 42: 30S ribosomal protein S7



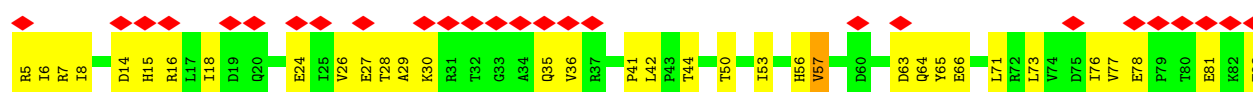
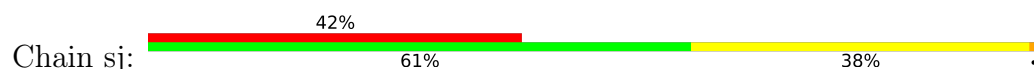
- Molecule 43: 30S ribosomal protein S8



- Molecule 44: Small ribosomal subunit protein uS9

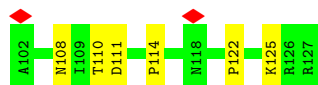
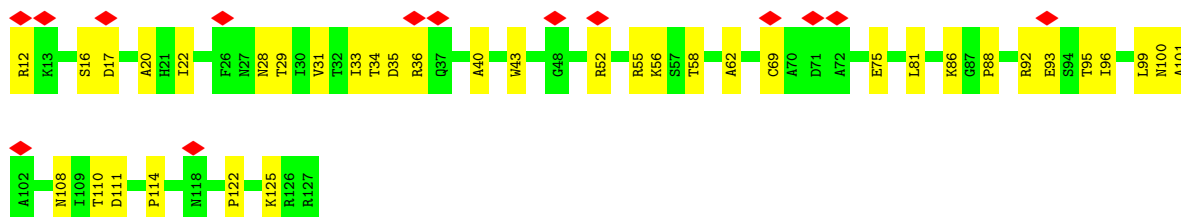


- Molecule 45: 30S ribosomal protein S10

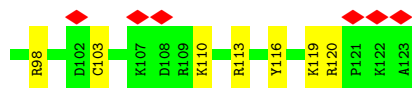
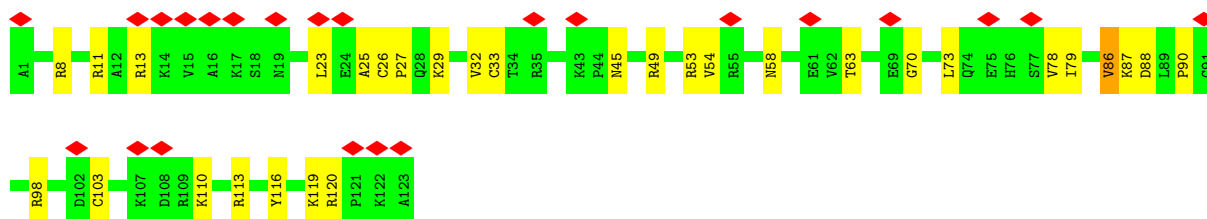
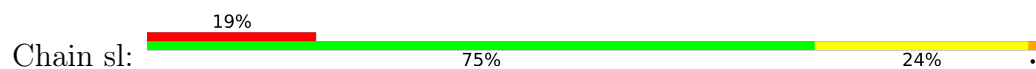




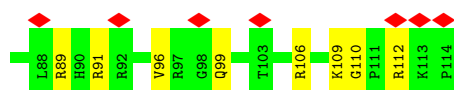
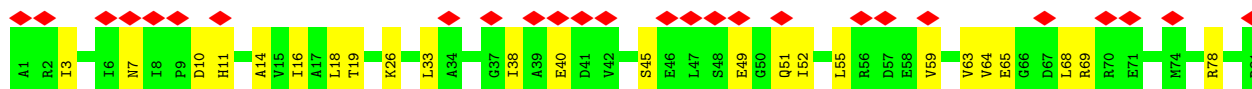
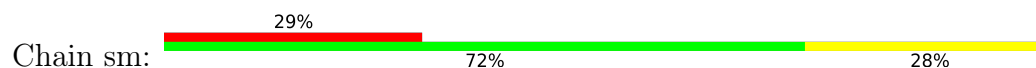
- Molecule 46: Small ribosomal subunit protein uS11



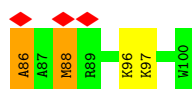
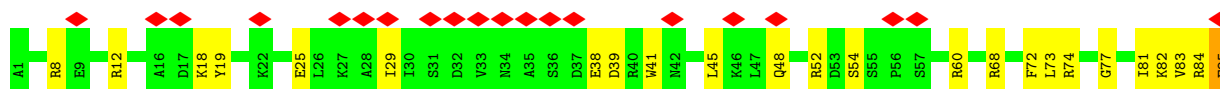
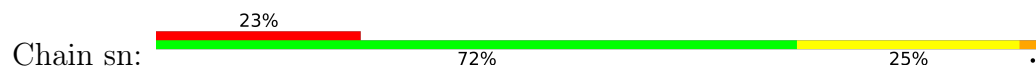
- Molecule 47: Small ribosomal subunit protein uS12



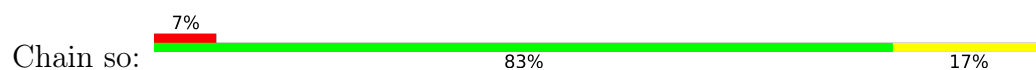
- Molecule 48: 30S ribosomal protein S13



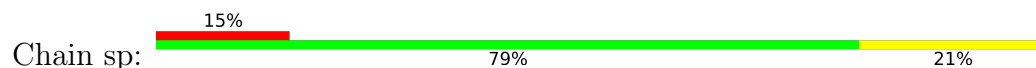
- Molecule 49: Small ribosomal subunit protein uS14



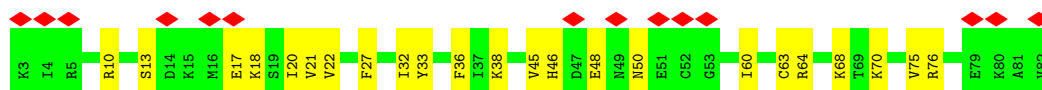
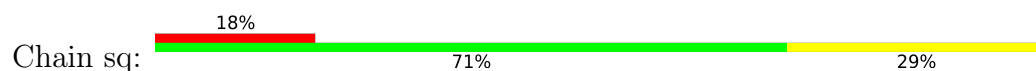
- Molecule 50: Small ribosomal subunit protein uS15



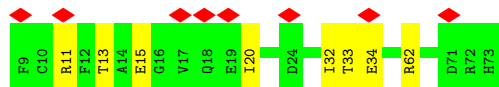
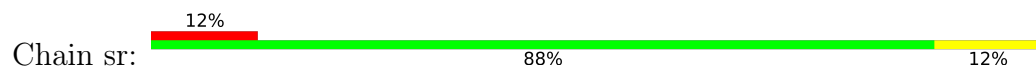
- Molecule 51: Small ribosomal subunit protein bS16



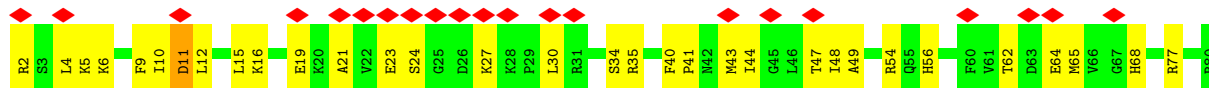
- Molecule 52: Small ribosomal subunit protein uS17



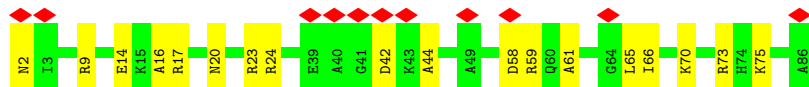
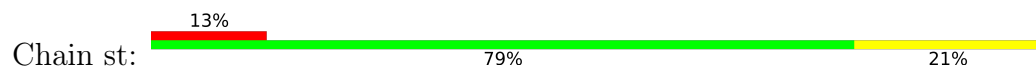
- Molecule 53: 30S ribosomal protein S18



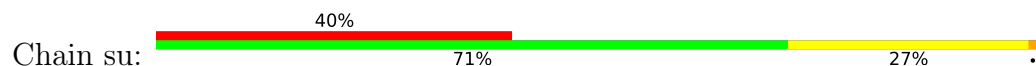
- Molecule 54: 30S ribosomal protein S19



- Molecule 55: 30S ribosomal protein S20



- Molecule 56: Small ribosomal subunit protein bS21





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	60175	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	64	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.569	Depositor
Minimum map value	-1.287	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.123	Depositor
Recommended contour level	0.45	Depositor
Map size (Å)	380.0, 380.0, 380.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.95, 0.95, 0.95	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PSU, H2U, MG, 4SU, MUM, NA, 4OC, FME, ATP

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	1	0.11	0/1361	0.39	0/1796
2	13	0.14	0/1152	0.29	0/1551
3	14	0.15	0/947	0.36	0/1268
4	15	0.15	0/1054	0.32	0/1403
5	16	0.16	0/1093	0.40	0/1460
6	17	0.16	0/973	0.35	0/1301
7	18	0.12	0/902	0.31	0/1209
8	19	0.15	0/929	0.30	0/1242
9	2	0.16	0/2121	0.34	0/2852
10	20	0.15	0/960	0.27	0/1278
11	21	0.17	0/829	0.47	0/1107
12	22	0.15	0/864	0.29	0/1156
13	23	0.14	0/744	0.32	0/994
14	24	0.12	0/787	0.30	0/1051
15	25	0.15	0/766	0.32	0/1025
16	27	0.16	0/589	0.29	0/779
17	28	0.14	0/635	0.30	0/848
18	29	0.13	0/510	0.27	0/677
19	3	0.15	0/1586	0.34	0/2134
20	30	0.14	0/453	0.26	0/605
21	31	0.12	0/531	0.32	0/709
22	32	0.15	0/450	0.33	0/599
23	33	0.12	0/416	0.29	0/554
24	34	0.15	0/380	0.35	0/498
25	35	0.16	0/513	0.31	0/676
26	36	0.16	0/303	0.31	0/397
27	4	0.15	0/1571	0.32	0/2113
28	5	0.14	0/1434	0.37	0/1926
29	6	0.15	0/1343	0.41	1/1816 (0.1%)
30	9	0.13	0/1122	0.43	0/1515
31	E	0.16	0/4441	0.38	1/5993 (0.0%)
32	M	0.10	0/219	0.22	0/339

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	R1	0.15	0/69797	0.29	0/108890
34	R2	0.13	0/2847	0.28	0/4440
35	R3	0.14	0/36782	0.31	0/57377
36	T	0.17	0/1716	0.29	0/2672
37	sb	0.14	0/1735	0.39	0/2338
38	sc	0.12	0/1651	0.30	0/2225
39	sd	0.12	0/1665	0.33	0/2227
40	se	0.16	0/1169	0.45	0/1573
41	sf	0.15	0/835	0.49	0/1128
42	sg	0.19	0/1195	0.34	0/1602
43	sh	0.14	0/989	0.33	0/1326
44	si	0.14	0/1034	0.44	0/1375
45	sj	0.14	0/796	0.38	0/1077
46	sk	0.15	0/885	0.34	0/1195
47	sl	0.14	0/969	0.42	0/1300
48	sm	0.15	0/892	0.38	0/1193
49	sn	0.16	0/817	0.48	1/1088 (0.1%)
50	so	0.12	0/722	0.27	0/964
51	sp	0.22	0/659	0.46	0/884
52	sq	0.16	0/657	0.38	0/881
53	sr	0.12	0/544	0.28	0/731
54	ss	0.13	0/652	0.36	0/877
55	st	0.12	0/671	0.30	0/888
56	su	0.72	2/598 (0.3%)	0.79	3/792 (0.4%)
All	All	0.16	2/162255 (0.0%)	0.32	6/241914 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
13	23	0	1
24	34	0	1
29	6	0	1
30	9	0	1
31	E	0	1
36	T	8	0
39	sd	0	1
40	se	0	1
41	sf	0	2
47	sl	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
49	sn	0	2
All	All	8	12

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
56	su	11	PRO	CG-CD	-14.40	1.01	1.50
56	su	11	PRO	N-CD	6.51	1.56	1.47

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	su	11	PRO	N-CD-CG	-12.71	84.14	103.20
56	su	11	PRO	CA-N-CD	-11.77	95.52	112.00
31	E	67	PRO	CA-N-CD	-10.54	97.25	112.00
56	su	11	PRO	CA-CB-CG	-7.97	89.36	104.50
29	6	11	PRO	CA-N-CD	-5.32	104.55	112.00

5 of 8 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
36	T	8	4SU	C2',C1'
36	T	20	H2U	C2'
36	T	32	4OC	C2'
36	T	54	MUM	C5,C4',C3'
36	T	55	PSU	C4'

5 of 12 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
13	23	37	ASP	Peptide
24	34	44	VAL	Peptide
29	6	12	ALA	Peptide
30	9	8	LYS	Peptide
31	E	16	PRO	Peptide

5.2 Too-close contacts

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	1353	0	1159	37	0
2	13	1129	0	1162	27	0
3	14	938	0	1012	21	0
4	15	1045	0	1117	24	0
5	16	1074	0	1157	24	0
6	17	960	0	1000	15	0
7	18	892	0	923	15	0
8	19	917	0	965	16	0
9	2	2082	0	2157	45	0
10	20	947	0	1022	13	0
11	21	816	0	839	23	0
12	22	857	0	922	18	0
13	23	738	0	807	15	0
14	24	779	0	834	7	0
15	25	753	0	780	22	0
16	27	582	0	599	20	0
17	28	625	0	655	14	0
18	29	509	0	543	8	0
19	3	1565	0	1616	35	0
20	30	449	0	491	11	0
21	31	522	0	524	14	0
22	32	444	0	461	12	0
23	33	409	0	440	10	0
24	34	377	0	418	6	0
25	35	504	0	574	12	0
26	36	302	0	343	9	0
27	4	1552	0	1619	26	0
28	5	1410	0	1447	39	0
29	6	1323	0	1374	32	0
30	9	1111	0	1148	30	0
31	E	4360	0	4336	102	0
32	M	195	0	99	0	0
33	R1	62318	0	31345	707	0
34	R2	2546	0	1292	34	0
35	R3	32850	0	16534	445	0
36	T	1639	0	831	28	0
37	sb	1704	0	1732	41	0
38	sc	1624	0	1699	26	0
39	sd	1643	0	1710	32	0
40	se	1156	0	1199	35	0
41	sf	817	0	808	21	0
42	sg	1181	0	1238	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
43	sh	979	0	1034	26	0
44	si	1022	0	1070	39	0
45	sj	786	0	828	24	0
46	sk	869	0	878	32	0
47	sl	955	0	1019	26	0
48	sm	883	0	944	20	0
49	sn	805	0	847	27	0
50	so	714	0	737	14	0
51	sp	649	0	666	14	0
52	sq	648	0	691	18	0
53	sr	535	0	552	5	0
54	ss	637	0	665	28	0
55	st	665	0	714	17	0
56	su	590	0	629	18	0
57	15	1	0	0	0	0
57	17	1	0	0	0	0
57	M	1	0	0	0	0
57	R1	204	0	0	0	0
57	R3	89	0	0	0	0
57	sn	1	0	0	0	0
57	sq	1	0	0	0	0
58	E	62	0	22	3	0
59	E	2	0	0	0	0
60	T	10	0	10	1	0
All	All	150106	0	102237	2142	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

The worst 5 of 2142 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:R3:1130:A:HO2'	35:R3:1131:G:H8	1.02	0.95
35:R3:359:G:HO2'	35:R3:360:G:H8	0.98	0.95
33:R1:132:G:H1	33:R1:146:A:H62	0.94	0.92
33:R1:2099:U:H3	33:R1:2190:G:H1	0.92	0.89
33:R1:1043:C:O2	33:R1:1112:G:N2	2.05	0.89

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1	218/220 (99%)	196 (90%)	22 (10%)	0	100	100
2	13	140/142 (99%)	138 (99%)	2 (1%)	0	100	100
3	14	120/122 (98%)	108 (90%)	12 (10%)	0	100	100
4	15	141/143 (99%)	130 (92%)	11 (8%)	0	100	100
5	16	134/136 (98%)	127 (95%)	7 (5%)	0	100	100
6	17	118/120 (98%)	109 (92%)	9 (8%)	0	100	100
7	18	114/116 (98%)	110 (96%)	4 (4%)	0	100	100
8	19	112/114 (98%)	101 (90%)	11 (10%)	0	100	100
9	2	269/271 (99%)	253 (94%)	16 (6%)	0	100	100
10	20	115/117 (98%)	115 (100%)	0	0	100	100
11	21	101/103 (98%)	88 (87%)	13 (13%)	0	100	100
12	22	108/110 (98%)	100 (93%)	8 (7%)	0	100	100
13	23	91/93 (98%)	82 (90%)	9 (10%)	0	100	100
14	24	100/102 (98%)	87 (87%)	13 (13%)	0	100	100
15	25	92/94 (98%)	89 (97%)	3 (3%)	0	100	100
16	27	74/76 (97%)	69 (93%)	5 (7%)	0	100	100
17	28	75/77 (97%)	74 (99%)	1 (1%)	0	100	100
18	29	61/63 (97%)	60 (98%)	1 (2%)	0	100	100
19	3	207/209 (99%)	195 (94%)	12 (6%)	0	100	100
20	30	56/58 (97%)	53 (95%)	3 (5%)	0	100	100
21	31	64/66 (97%)	55 (86%)	9 (14%)	0	100	100
22	32	54/56 (96%)	51 (94%)	3 (6%)	0	100	100
23	33	48/50 (96%)	48 (100%)	0	0	100	100
24	34	44/46 (96%)	40 (91%)	4 (9%)	0	100	100
25	35	62/64 (97%)	59 (95%)	3 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	36	36/38 (95%)	34 (94%)	2 (6%)	0	100	100
27	4	199/201 (99%)	188 (94%)	11 (6%)	0	100	100
28	5	175/177 (99%)	161 (92%)	14 (8%)	0	100	100
29	6	174/176 (99%)	159 (91%)	15 (9%)	0	100	100
30	9	147/149 (99%)	135 (92%)	10 (7%)	2 (1%)	9	39
31	E	550/552 (100%)	508 (92%)	41 (8%)	1 (0%)	43	73
37	sb	216/218 (99%)	194 (90%)	21 (10%)	1 (0%)	24	59
38	sc	204/206 (99%)	197 (97%)	7 (3%)	0	100	100
39	sd	203/205 (99%)	185 (91%)	18 (9%)	0	100	100
40	se	155/157 (99%)	135 (87%)	18 (12%)	2 (1%)	9	40
41	sf	98/100 (98%)	83 (85%)	14 (14%)	1 (1%)	12	45
42	sg	149/151 (99%)	142 (95%)	7 (5%)	0	100	100
43	sh	127/129 (98%)	121 (95%)	6 (5%)	0	100	100
44	si	125/127 (98%)	102 (82%)	22 (18%)	1 (1%)	16	50
45	sj	96/98 (98%)	88 (92%)	7 (7%)	1 (1%)	12	45
46	sk	114/116 (98%)	103 (90%)	11 (10%)	0	100	100
47	sl	121/123 (98%)	92 (76%)	29 (24%)	0	100	100
48	sm	112/114 (98%)	99 (88%)	13 (12%)	0	100	100
49	sn	98/100 (98%)	78 (80%)	20 (20%)	0	100	100
50	so	86/88 (98%)	83 (96%)	3 (4%)	0	100	100
51	sp	80/82 (98%)	71 (89%)	9 (11%)	0	100	100
52	sq	78/80 (98%)	67 (86%)	11 (14%)	0	100	100
53	sr	63/65 (97%)	61 (97%)	2 (3%)	0	100	100
54	ss	77/79 (98%)	71 (92%)	6 (8%)	0	100	100
55	st	83/85 (98%)	82 (99%)	1 (1%)	0	100	100
56	su	68/70 (97%)	64 (94%)	3 (4%)	1 (2%)	8	37
All	All	6352/6454 (98%)	5840 (92%)	502 (8%)	10 (0%)	44	73

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
31	E	17	PRO
30	9	8	LYS

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Mol	Chain	Res	Type
44	si	91	GLU
37	sb	18	GLN
40	se	122	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	106/171 (62%)	105 (99%)	1 (1%)	70	81
2	13	116/116 (100%)	116 (100%)	0	100	100
3	14	103/103 (100%)	101 (98%)	2 (2%)	50	73
4	15	102/102 (100%)	101 (99%)	1 (1%)	68	80
5	16	109/109 (100%)	108 (99%)	1 (1%)	70	81
6	17	100/100 (100%)	100 (100%)	0	100	100
7	18	86/86 (100%)	86 (100%)	0	100	100
8	19	99/99 (100%)	98 (99%)	1 (1%)	68	80
9	2	216/216 (100%)	216 (100%)	0	100	100
10	20	89/89 (100%)	89 (100%)	0	100	100
11	21	84/84 (100%)	83 (99%)	1 (1%)	63	79
12	22	93/93 (100%)	93 (100%)	0	100	100
13	23	80/80 (100%)	80 (100%)	0	100	100
14	24	83/83 (100%)	83 (100%)	0	100	100
15	25	78/78 (100%)	78 (100%)	0	100	100
16	27	58/58 (100%)	58 (100%)	0	100	100
17	28	67/67 (100%)	67 (100%)	0	100	100
18	29	55/55 (100%)	55 (100%)	0	100	100
19	3	164/164 (100%)	163 (99%)	1 (1%)	78	84
20	30	48/48 (100%)	48 (100%)	0	100	100
21	31	59/59 (100%)	59 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
22	32	47/47 (100%)	47 (100%)	0	100	100
23	33	45/45 (100%)	44 (98%)	1 (2%)	45	71
24	34	38/38 (100%)	38 (100%)	0	100	100
25	35	51/51 (100%)	51 (100%)	0	100	100
26	36	34/34 (100%)	33 (97%)	1 (3%)	37	67
27	4	165/165 (100%)	165 (100%)	0	100	100
28	5	148/148 (100%)	147 (99%)	1 (1%)	76	83
29	6	137/137 (100%)	137 (100%)	0	100	100
30	9	114/114 (100%)	113 (99%)	1 (1%)	70	81
31	E	461/464 (99%)	459 (100%)	2 (0%)	84	86
37	sb	180/180 (100%)	180 (100%)	0	100	100
38	sc	170/170 (100%)	170 (100%)	0	100	100
39	sd	172/172 (100%)	171 (99%)	1 (1%)	78	84
40	se	119/119 (100%)	119 (100%)	0	100	100
41	sf	87/87 (100%)	83 (95%)	4 (5%)	24	58
42	sg	124/124 (100%)	124 (100%)	0	100	100
43	sh	104/104 (100%)	104 (100%)	0	100	100
44	si	105/105 (100%)	105 (100%)	0	100	100
45	sj	86/86 (100%)	84 (98%)	2 (2%)	44	70
46	sk	89/89 (100%)	89 (100%)	0	100	100
47	sl	103/103 (100%)	103 (100%)	0	100	100
48	sm	92/92 (100%)	92 (100%)	0	100	100
49	sn	83/83 (100%)	83 (100%)	0	100	100
50	so	76/76 (100%)	76 (100%)	0	100	100
51	sp	65/65 (100%)	64 (98%)	1 (2%)	57	76
52	sq	74/74 (100%)	74 (100%)	0	100	100
53	sr	56/56 (100%)	55 (98%)	1 (2%)	51	74
54	ss	70/70 (100%)	69 (99%)	1 (1%)	59	77
55	st	65/65 (100%)	65 (100%)	0	100	100
56	su	60/60 (100%)	60 (100%)	0	100	100
All	All	5215/5283 (99%)	5191 (100%)	24 (0%)	78	85

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
39	sd	142	VAL
41	sf	81	ASN
41	sf	53	LYS
41	sf	94	HIS
11	21	26	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 56 such sidechains are listed below:

Mol	Chain	Res	Type
28	5	126	ASN
55	st	81	GLN
31	E	133	GLN
55	st	19	HIS
44	si	36	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
32	M	8/9 (88%)	0	0
33	R1	2902/2903 (99%)	570 (19%)	18 (0%)
34	R2	118/119 (99%)	19 (16%)	0
35	R3	1529/1531 (99%)	399 (26%)	23 (1%)
36	T	76/77 (98%)	20 (26%)	3 (3%)
All	All	4633/4639 (99%)	1008 (21%)	44 (0%)

5 of 1008 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
33	R1	2	G
33	R1	5	A
33	R1	10	A
33	R1	15	G
33	R1	34	U

5 of 44 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
35	R3	500	G
35	R3	837	U

Continued on next page...

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Mol	Chain	Res	Type
35	R3	561	U
35	R3	672	U
35	R3	1149	C

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

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
36	H2U	T	20	36	18,21,22	4.50	5 (27%)	19,30,33	4.25	6 (31%)
36	MUM	T	54	36	18,22,22	3.27	6 (33%)	19,32,32	2.11	5 (26%)
36	4SU	T	8	36	18,21,22	3.72	7 (38%)	25,30,33	2.41	6 (24%)
36	4OC	T	32	36	20,23,24	2.53	5 (25%)	25,32,35	0.94	1 (4%)
36	PSU	T	55	36	18,21,22	2.30	8 (44%)	21,30,33	1.85	3 (14%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
36	H2U	T	20	36	1/1/8/9	6/7/38/39	0/2/2/2
36	MUM	T	54	36	3/3/9/10	4/7/41/41	0/2/2/2
36	4SU	T	8	36	2/2/5/5	2/7/25/26	0/2/2/2
36	4OC	T	32	36	1/1/5/6	1/9/29/30	0/2/2/2
36	PSU	T	55	36	1/1/5/5	3/7/25/26	0/2/2/2

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
36	T	54	MUM	C6-N1	-11.96	1.32	1.46

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
36	T	8	4SU	O2-C2	10.81	1.42	1.23
36	T	20	H2U	C2-N1	10.52	1.50	1.35
36	T	20	H2U	O4-C4	10.12	1.43	1.23
36	T	32	4OC	O2-C2	8.92	1.40	1.23

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
36	T	20	H2U	O2-C2-N1	-12.74	107.78	123.10
36	T	20	H2U	O4-C4-N3	-7.30	109.04	120.30
36	T	8	4SU	C4-N3-C2	-7.25	120.36	127.31
36	T	20	H2U	O2-C2-N3	-6.81	108.93	121.49
36	T	20	H2U	O4-C4-C5	-6.39	109.11	122.20

5 of 8 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
36	T	20	H2U	C2'
36	T	32	4OC	C2'
36	T	54	MUM	C5
36	T	54	MUM	C4'
36	T	54	MUM	C3'

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
36	T	20	H2U	O4'-C4'-C5'-O5'
36	T	20	H2U	O4'-C1'-N1-C2
36	T	20	H2U	O4'-C1'-N1-C6
36	T	54	MUM	O4'-C4'-C5'-O5'
36	T	54	MUM	C3'-C4'-C5'-O5'

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
36	T	20	H2U	1	0
36	T	8	4SU	1	0
36	T	32	4OC	1	0
36	T	55	PSU	2	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 303 ligands modelled in this entry, 300 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
60	FME	T	101	36	8,9,10	0.98	0	8,9,11	0.98	0
58	ATP	E	602	59	32,33,33	0.53	0	48,52,52	0.32	0
58	ATP	E	601	-	32,33,33	0.40	0	48,52,52	0.40	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
60	FME	T	101	36	-	1/7/9/11	-
58	ATP	E	602	59	-	2/22/38/38	0/3/3/3
58	ATP	E	601	-	-	9/22/38/38	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

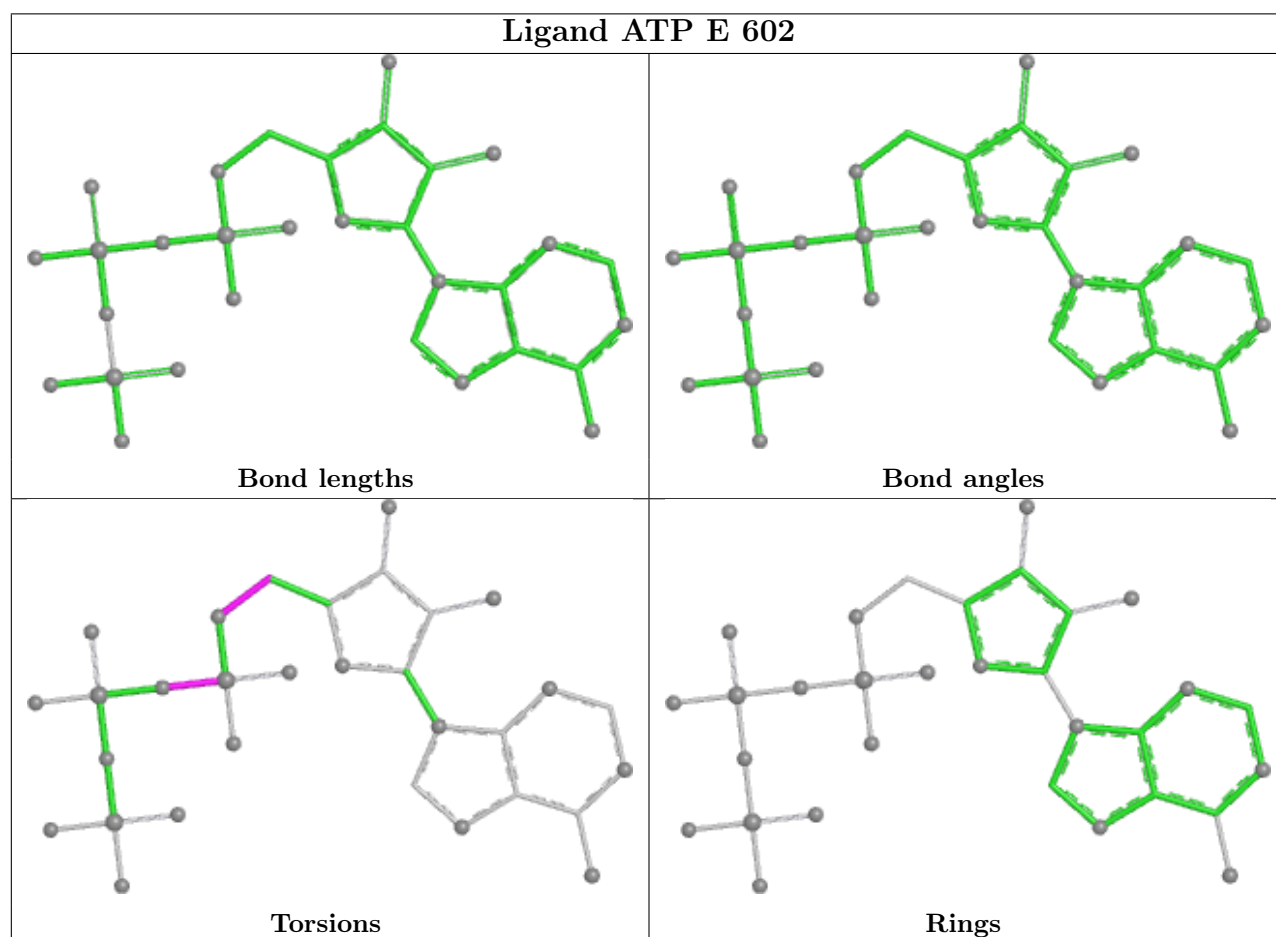
Mol	Chain	Res	Type	Atoms
58	E	601	ATP	C5'-O5'-PA-O1A
58	E	601	ATP	C5'-O5'-PA-O2A
58	E	601	ATP	C5'-O5'-PA-O3A
58	E	601	ATP	O4'-C4'-C5'-O5'
58	E	601	ATP	C3'-C4'-C5'-O5'

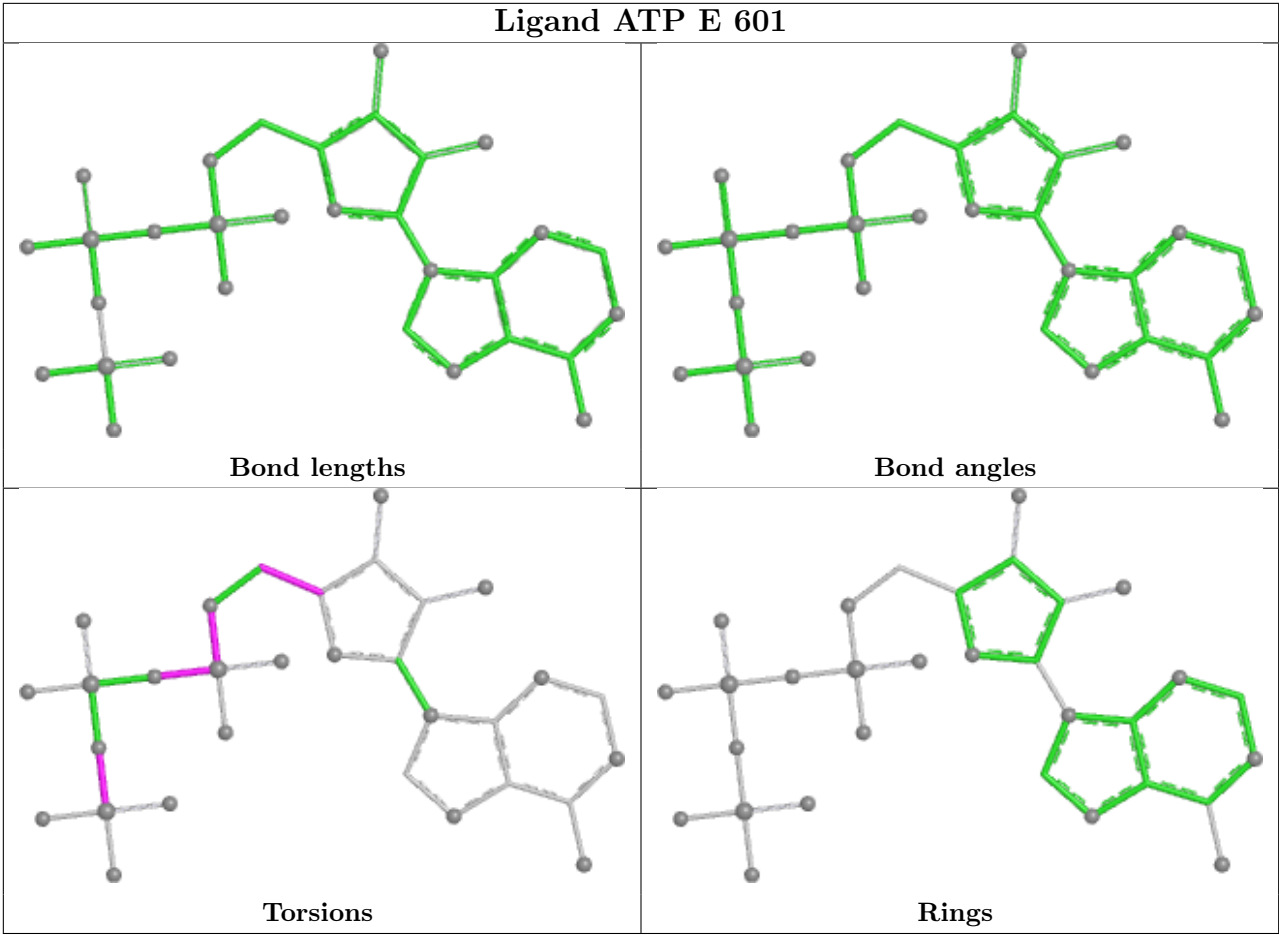
There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
60	T	101	FME	1	0
58	E	602	ATP	2	0
58	E	601	ATP	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
35	R3	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	R3	210:C	O3'	211:G	P	5.77
1	R3	460:A	O3'	461:A	P	5.31

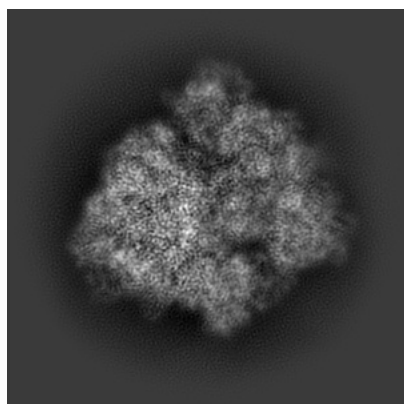
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29398. These allow visual inspection of the internal detail of the map and identification of artifacts.

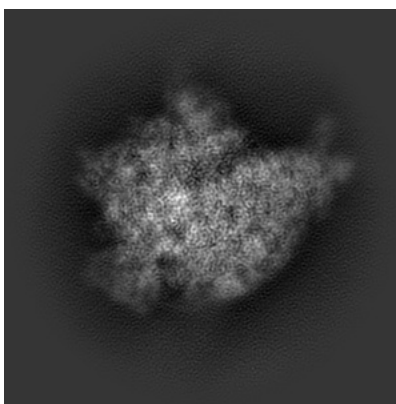
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

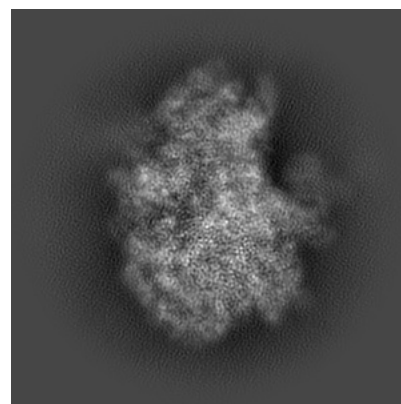
6.1.1 Primary map



X

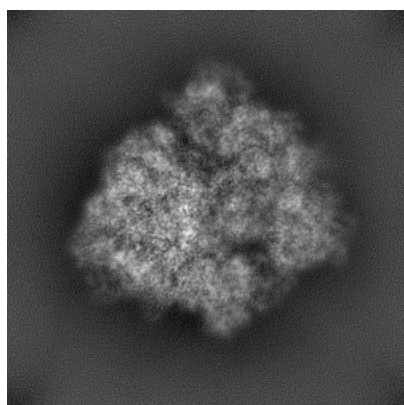


Y

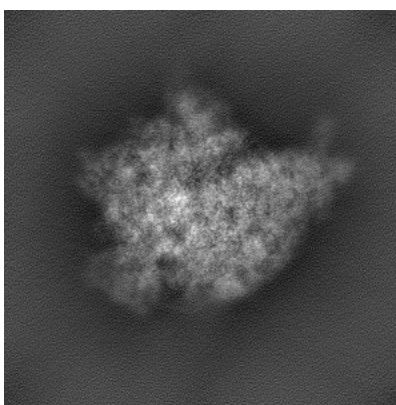


Z

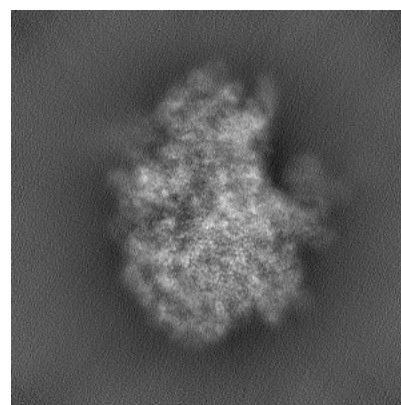
6.1.2 Raw map



X



Y

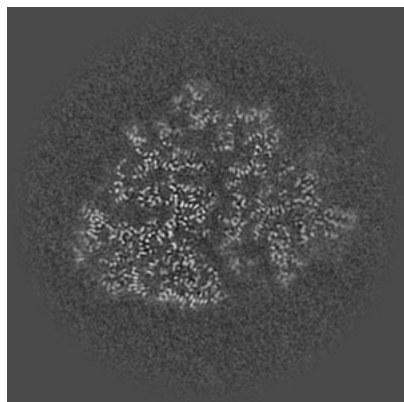


Z

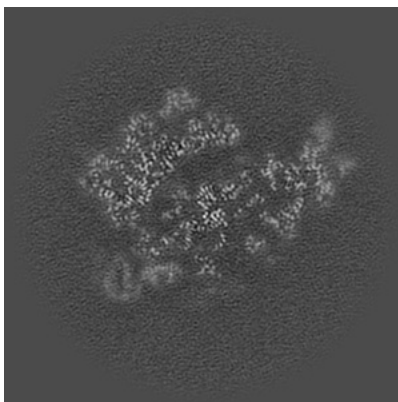
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

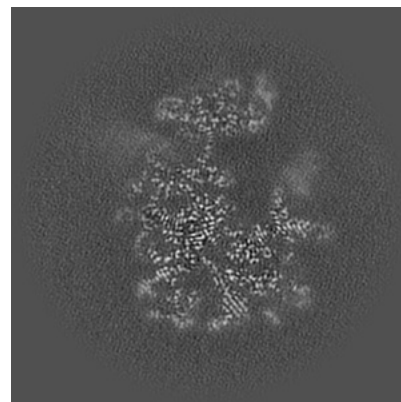
6.2.1 Primary map



X Index: 200

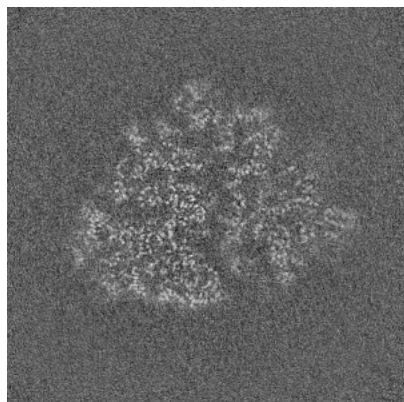


Y Index: 200

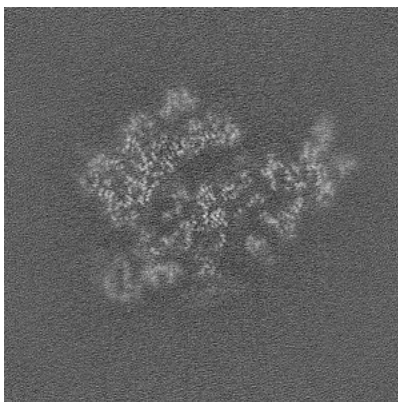


Z Index: 200

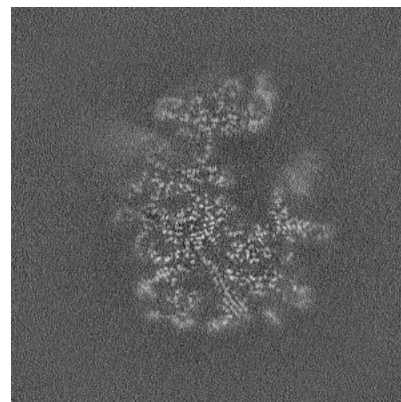
6.2.2 Raw map



X Index: 200



Y Index: 200

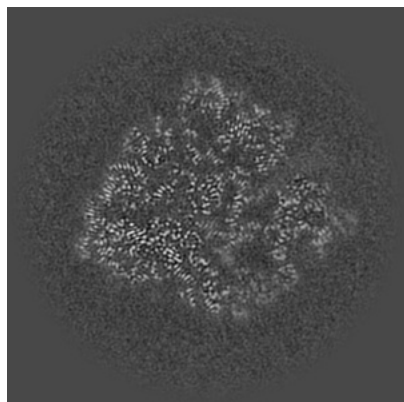


Z Index: 200

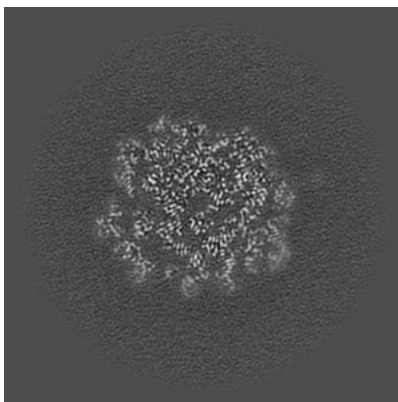
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

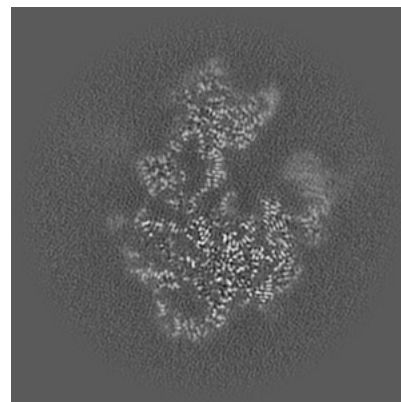
6.3.1 Primary map



X Index: 210

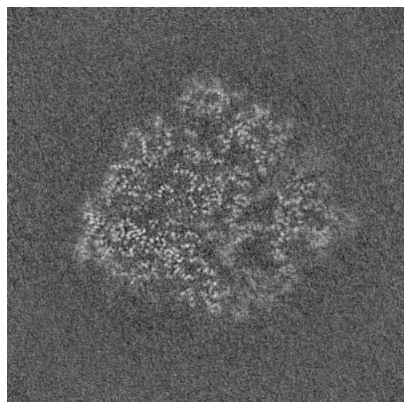


Y Index: 152

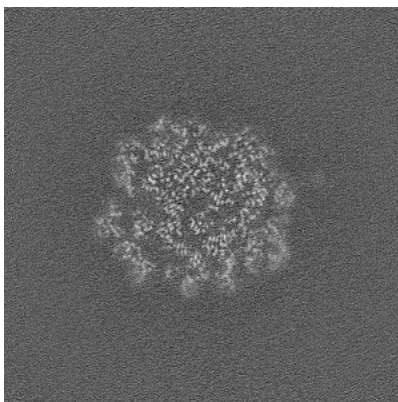


Z Index: 179

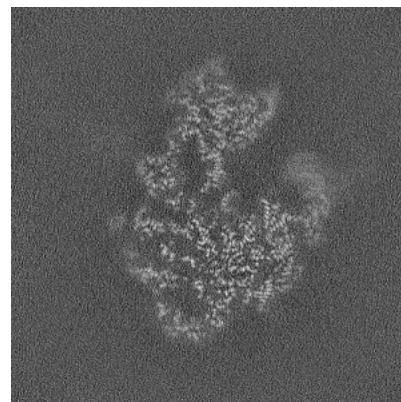
6.3.2 Raw map



X Index: 209



Y Index: 152

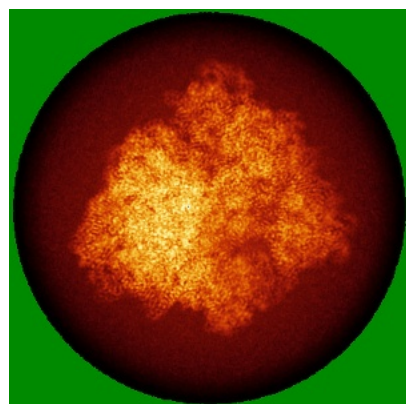


Z Index: 178

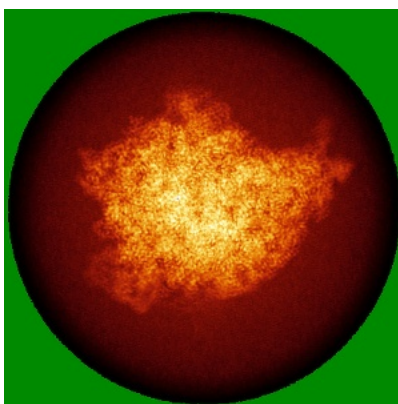
The images above show the largest variance slices of the map in three orthogonal directions.

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

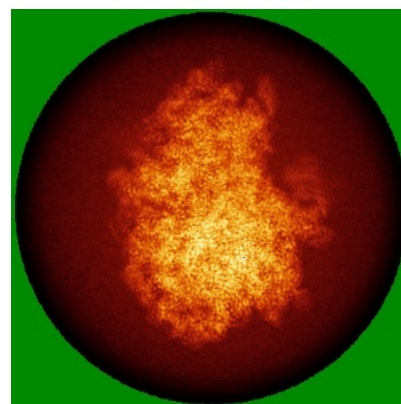
6.4.1 Primary map



X

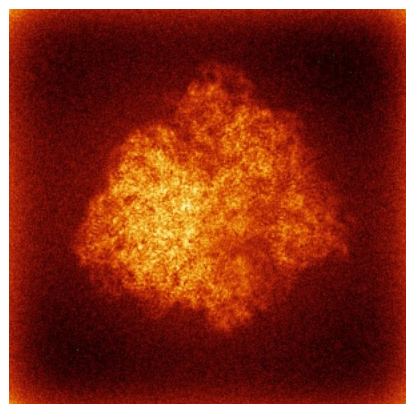


Y

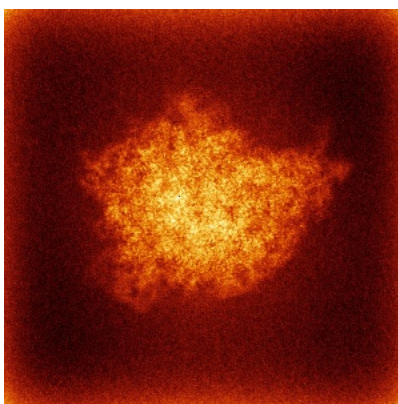


Z

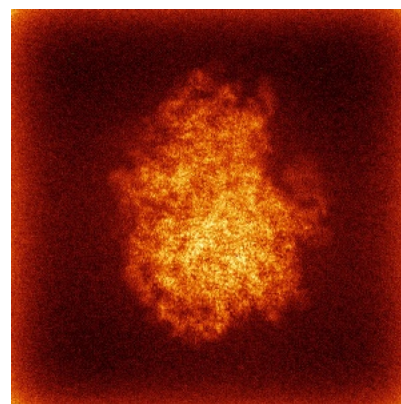
6.4.2 Raw map



X



Y

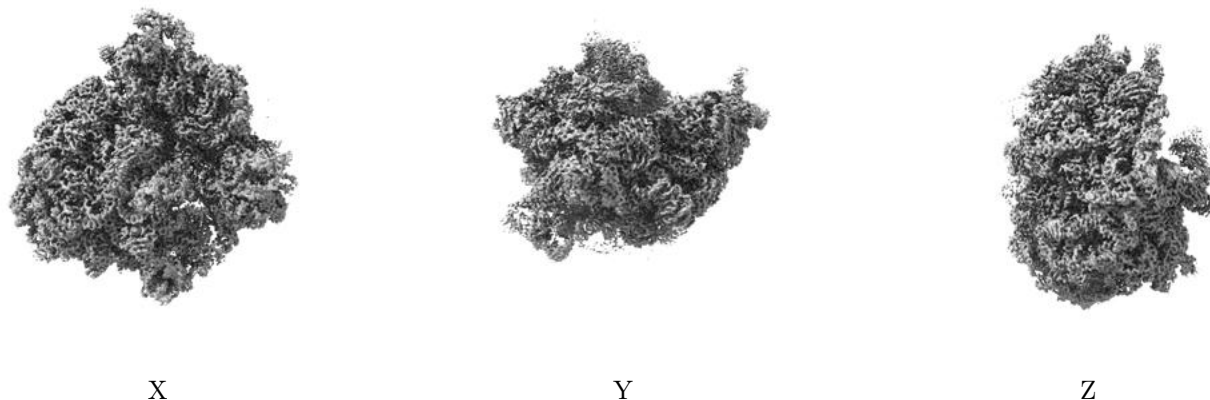


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.45. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

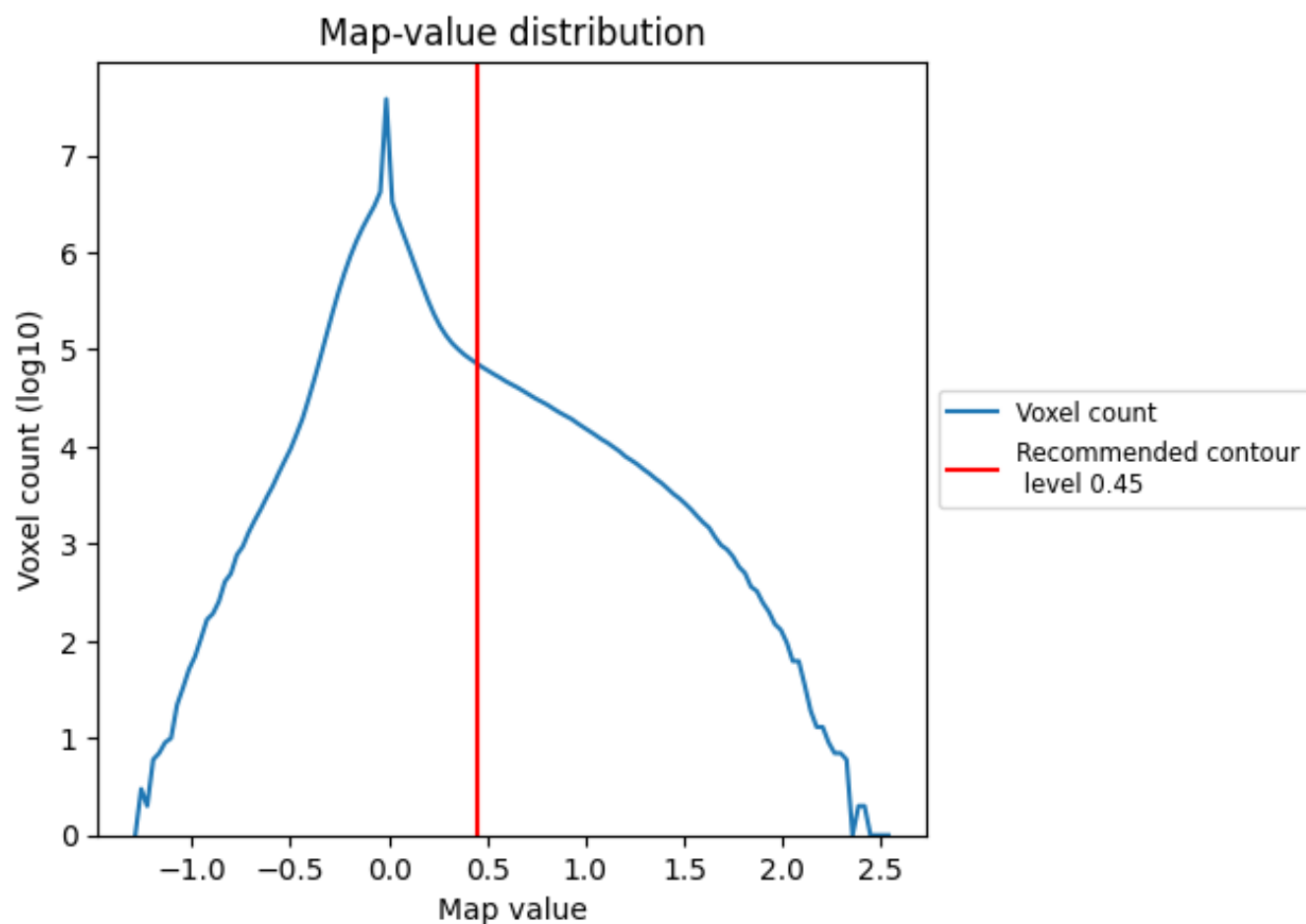
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

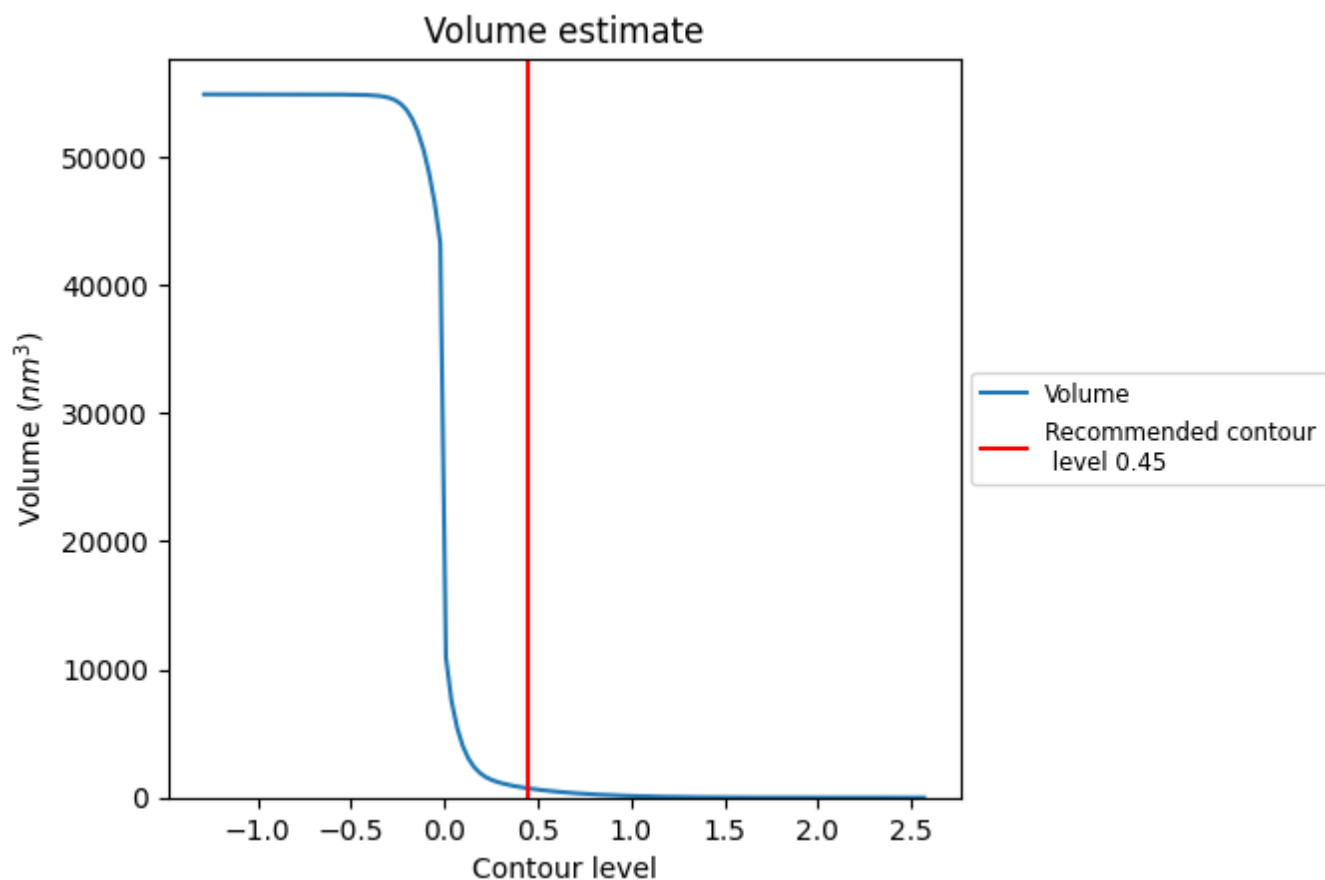
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

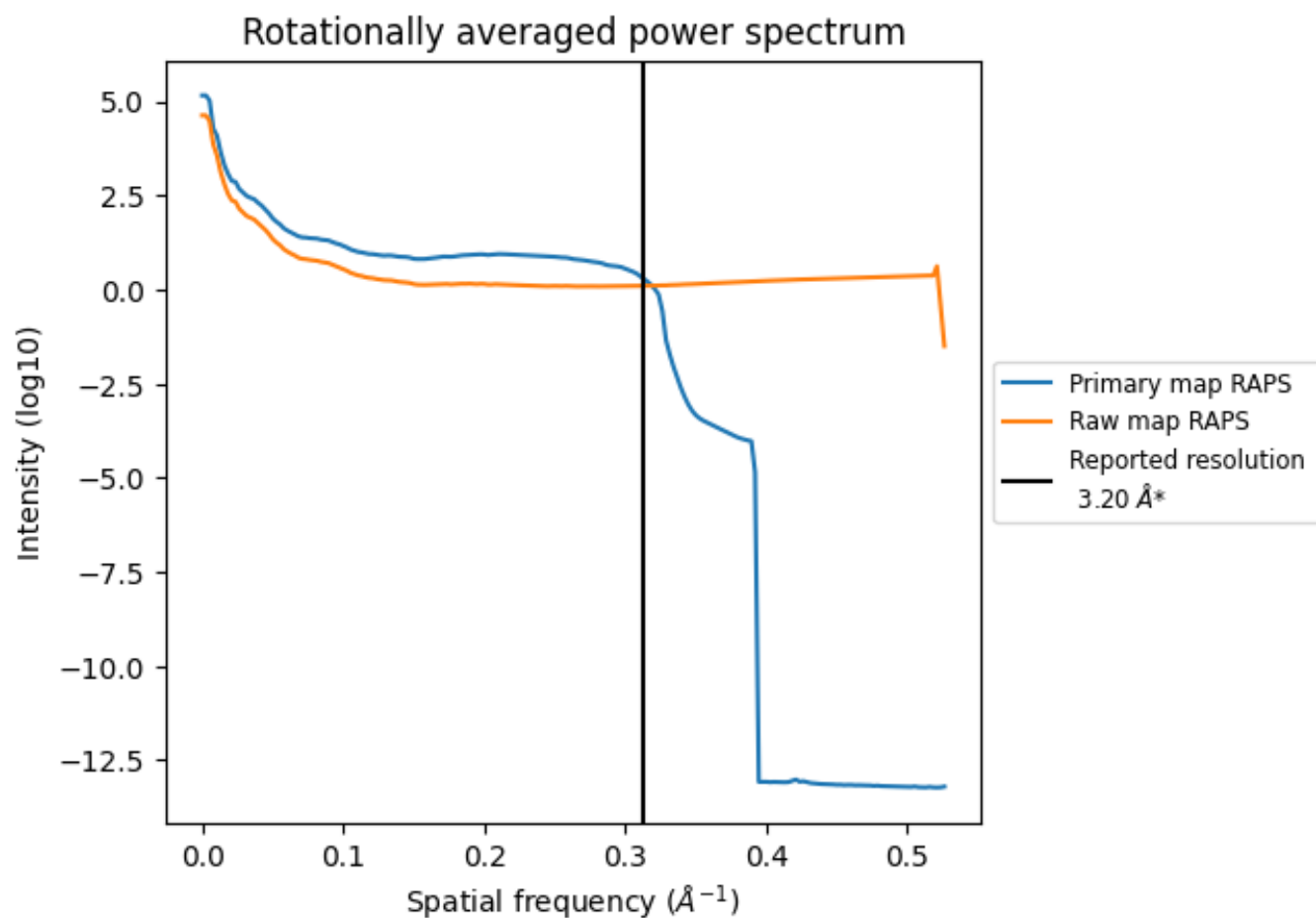
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 718 nm³; this corresponds to an approximate mass of 648 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

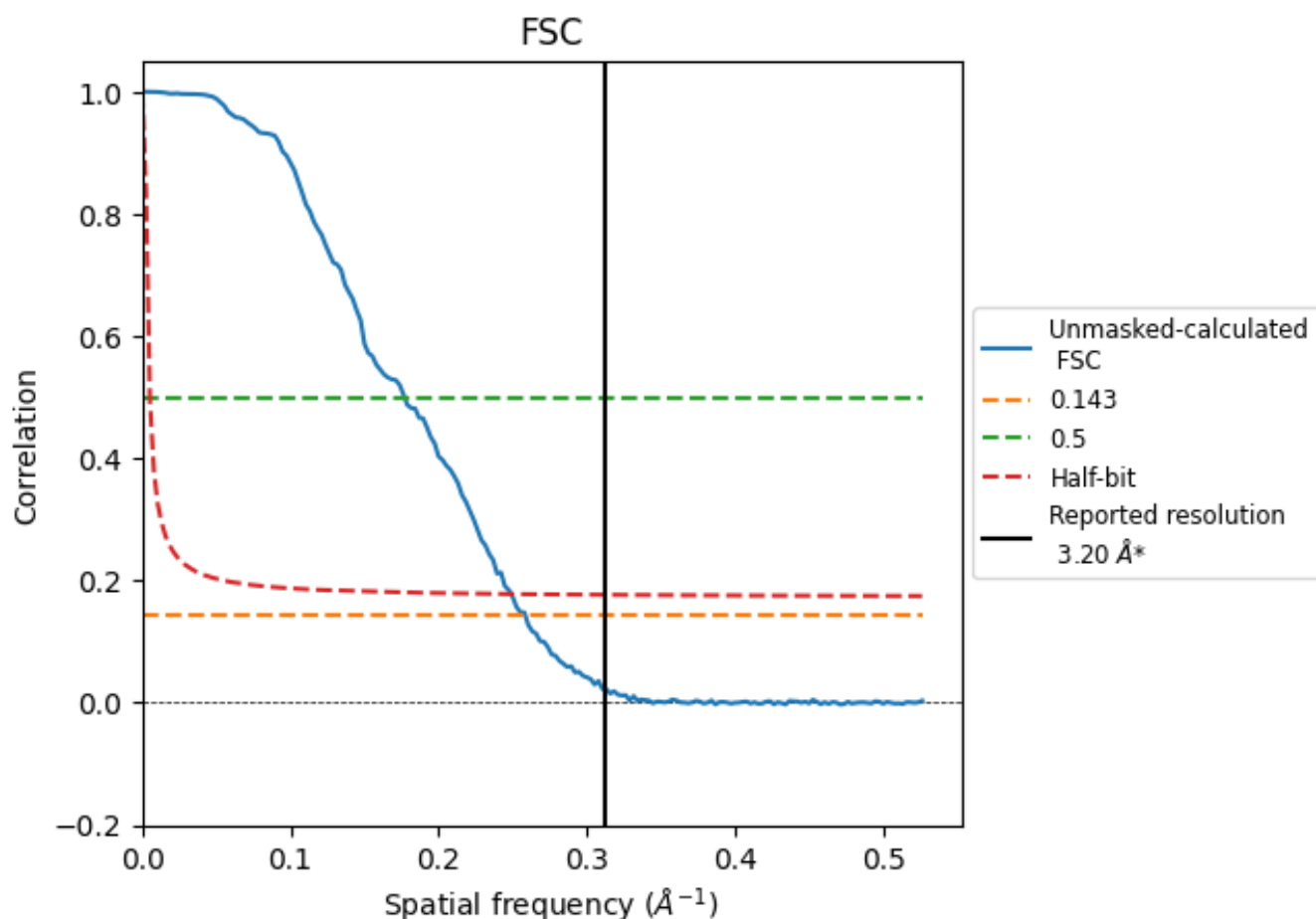


*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.87	5.65	4.02

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.87 differs from the reported value 3.2 by more than 10 %

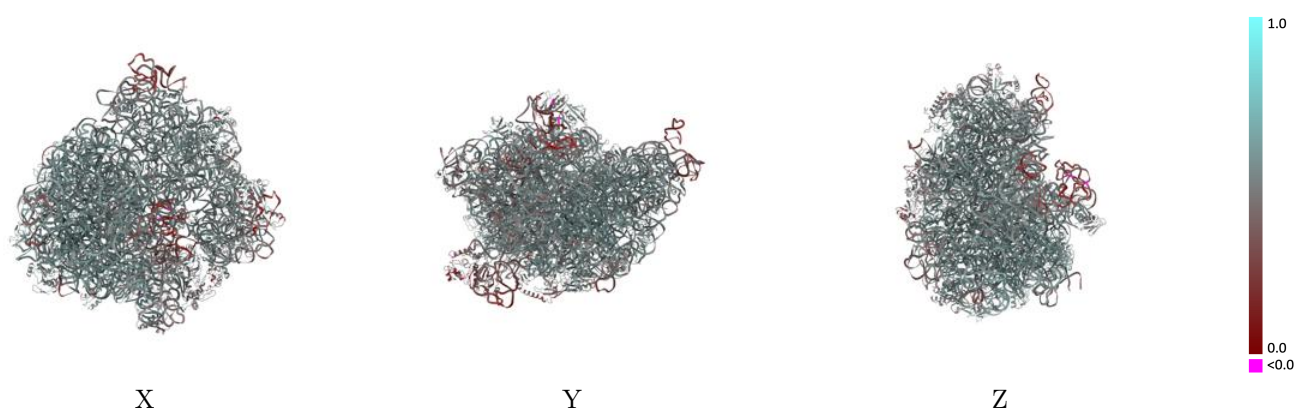
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29398 and PDB model 9NLE. Per-residue inclusion information can be found in section 3 on page 17.

9.1 Map-model overlay [i](#)

This section was not generated.

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

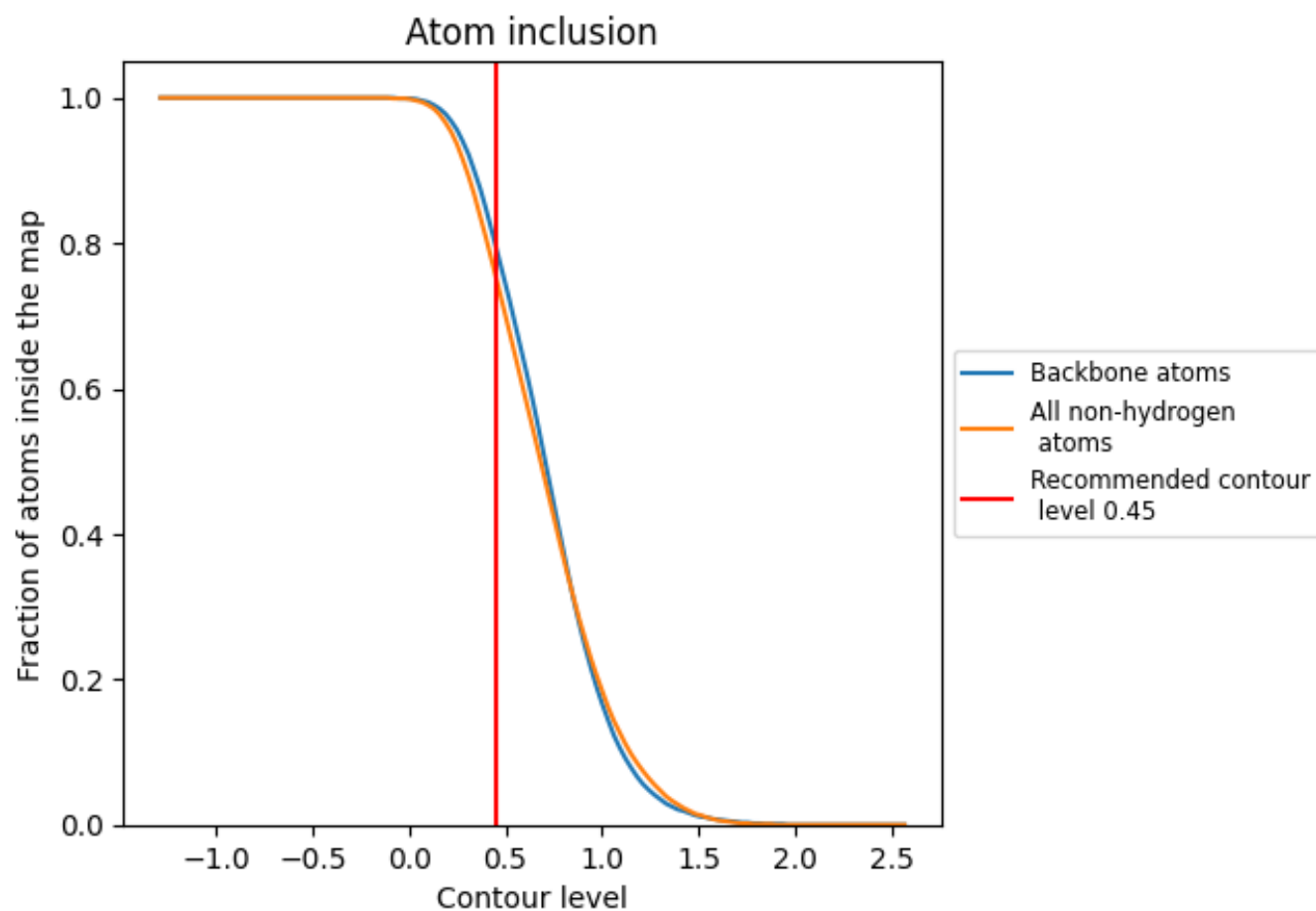


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 76% 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 (0.45) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7550	 0.5270
1	 0.1410	 0.3250
13	 0.7490	 0.5720
14	 0.6870	 0.5630
15	 0.7460	 0.5630
16	 0.7000	 0.5620
17	 0.7800	 0.5740
18	 0.6580	 0.5300
19	 0.6820	 0.5680
2	 0.7370	 0.5840
20	 0.8010	 0.5790
21	 0.7190	 0.5520
22	 0.6930	 0.5630
23	 0.6450	 0.5400
24	 0.6190	 0.5280
25	 0.6650	 0.5350
27	 0.7350	 0.5710
28	 0.7140	 0.5730
29	 0.5940	 0.5020
3	 0.7350	 0.5670
30	 0.7190	 0.5700
31	 0.2290	 0.4060
32	 0.7220	 0.5690
33	 0.6630	 0.5600
34	 0.7720	 0.5890
35	 0.7760	 0.5950
36	 0.7190	 0.5650
4	 0.6700	 0.5430
5	 0.4620	 0.4850
6	 0.5400	 0.4890
9	 0.2010	 0.4060
E	 0.5150	 0.5040
M	 0.5920	 0.5400
R1	 0.8420	 0.5360
R2	 0.8050	 0.4950



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Chain	Atom inclusion	Q-score
R3	 0.8140	 0.5180
T	 0.6180	 0.4590
sb	 0.5300	 0.4760
sc	 0.5870	 0.5330
sd	 0.6000	 0.5170
se	 0.6520	 0.5380
sf	 0.6070	 0.4990
sg	 0.5510	 0.5170
sh	 0.6650	 0.5520
si	 0.6040	 0.5060
sj	 0.4780	 0.4800
sk	 0.6500	 0.5510
sl	 0.6080	 0.5380
sm	 0.5600	 0.5070
sn	 0.5900	 0.4860
so	 0.6620	 0.5510
sp	 0.6540	 0.5300
sq	 0.6080	 0.5350
sr	 0.6610	 0.5360
ss	 0.5470	 0.4920
st	 0.6490	 0.5380
su	 0.4790	 0.5150