



Full wwPDB EM Validation Report ⓘ

Jan 30, 2025 – 02:54 PM JST

PDB ID : 8HL1
EMDB ID : EMD-34866
Title : Cryo-EM Structures and Translocation Mechanism of Crenarchaeota Ribosome
Authors : Wang, Y.H.; Zhou, J.
Deposited on : 2022-11-28
Resolution : 3.93 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

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.dev113
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.40

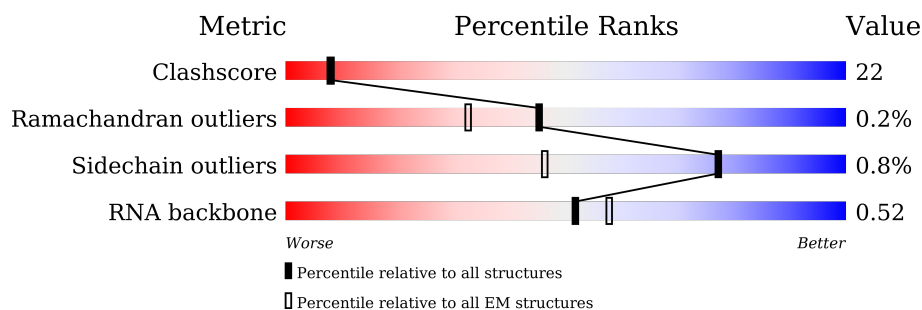
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.93 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

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	A16S	1494	 72% 27% .
2	A23S	3022	 76% 23% ..
3	AEFG	729	 14% 97% ..
4	A5S	122	 49% 39% 11%
5	AL2P	234	 12% 100%
6	AL3P	339	 11% 100%
7	AL4P	251	 6% 99% .

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Mol	Chain	Length	Quality of chain
8	AL5P	168	12% 99% .
9	AL6P	181	8% 100%
10	ALX0	76	13% 100%
11	L10E	164	11% 100%
12	L13P	140	11% 98% .
13	L141	86	8% 100%
13	L142	86	9% 100%
14	L14P	134	16% 99% .
15	L15E	169	6% 99% .
16	L18E	112	. 98% .
17	L18P	193	9% 99% .
18	L19E	144	. 100%
19	L22P	150	7% 100%
20	L23P	81	9% 100%
21	L24E	54	100%
22	L24P	122	9% 98% .
23	L29P	63	11% 100%
24	L30E	94	6% 99% .
25	L30P	155	13% 100%
26	L31E	75	5% 100%
27	L32E	123	18% 98% .
28	L34E	77	26% 96% .
29	L37A	65	8% 98% .
30	L37E	54	11% 100%
31	L39E	49	14% 98% .

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Mol	Chain	Length	Quality of chain
32	L40E	55	 35% 96%
33	L44E	92	 100%
34	L7A1	123	 7% 100%
34	L7A2	123	 28% 100%
34	SL7A	123	 89% 99%
35	L15P	144	 65% 35%
36	L21E	97	 6% 100%
37	L45A	101	 20% 95% 5%
38	L46A	70	 29% 99%
39	L47A	80	 95% 99%
40	AL1P	216	 63% 99%
41	AS3P	201	 30% 100%
42	AS7P	193	 58% 98%
43	S10P	100	 47% 97%
44	S13P	147	 56% 100%
45	S14P	52	 29% 100%
46	S17E	62	 42% 100%
47	S19E	150	 55% 99%
48	S19P	115	 36% 99%
49	AS9P	136	 40% 100%
50	S28E	63	 48% 100%
51	S27A	54	 30% 98%
52	AS2P	196	 49% 99%
53	AS4E	240	 11% 98%
54	AS4P	166	 8% 99%

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Mol	Chain	Length	Quality of chain
55	AS5P	204	<div> <div>8%</div> <div>100%</div> </div>
56	AS6E	105	<div> <div>22%</div> <div>98%</div> </div>
57	AS8E	126	<div> <div>17%</div> <div>99%</div> </div>
58	AS8P	130	<div> <div>7%</div> <div>99%</div> </div>
59	S11P	128	<div> <div>15%</div> <div>100%</div> </div>
60	S12P	143	<div> <div>15%</div> <div>99%</div> </div>
61	S15P	149	<div> <div>9%</div> <div>99%</div> </div>
62	S17P	111	<div> <div>11%</div> <div>98%</div> </div>
63	S24E	96	<div> <div>•</div> <div>99%</div> </div>
64	S27E	59	<div> <div>•</div> <div>100%</div> </div>
65	S3AE	189	<div> <div>12%</div> <div>97%</div> </div>
66	AETN	76	<div> <div>7%</div> <div>53%</div> <div>42%</div> <div>5%</div> </div>
66	APTN	76	<div> <div>•</div> <div>71%</div> <div>25%</div> <div>•</div> </div>
67	AMRN	13	<div> <div>23%</div> <div>46%</div> <div>54%</div> </div>
68	APTP	6	<div> <div>83%</div> <div>100%</div> </div>

2 Entry composition

There are 70 unique types of molecules in this entry. The entry contains 173502 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 16s rRNA (1493-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A16S	1493	Total	C	N	O	P	0	0
			32063	14279	5930	10361	1493		

- Molecule 2 is a RNA chain called 23s rRNA (3000-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A23S	2995	Total	C	N	O	P	0	0
			64339	28665	11911	20768	2995		

- Molecule 3 is a protein called Elongation factor 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	AEFG	725	Total	C	N	O	S	0	0
			5677	3609	984	1067	17		

- Molecule 4 is a RNA chain called 5s rRNA (122-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	A5S	122	Total	C	N	O	P	0	0
			2609	1163	476	849	121		

- Molecule 5 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	AL2P	234	Total	C	N	O	S	0	0
			1754	1101	344	307	2		

- Molecule 6 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	AL3P	339	Total	C	N	O	S	0	0
			2695	1730	484	477	4		

- Molecule 7 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	AL4P	251	Total	C	N	O	S	0	0
			1926	1223	356	345	2		

- Molecule 8 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	AL5P	168	Total	C	N	O	S	0	0
			1343	854	253	232	4		

- Molecule 9 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	AL6P	181	Total	C	N	O	S	0	0
			1431	920	246	264	1		

- Molecule 10 is a protein called 50S ribosomal protein L18Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	ALX0	76	Total	C	N	O	S	0	0
			629	403	110	115	1		

- Molecule 11 is a protein called 50S ribosomal protein L10e.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	L10E	164	Total	C	N	O	S	0	0
			1310	837	239	227	7		

- Molecule 12 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L13P	140	Total	C	N	O	S	0	0
			1109	707	208	190	4		

- Molecule 13 is a protein called 50S ribosomal protein L14e.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	L141	86	Total	C	N	O	S	0	0
			669	417	123	127	2		
13	L142	86	Total	C	N	O	S	0	0
			669	417	123	127	2		

- Molecule 14 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	L14P	134	Total	C	N	O	S	0	0
			1034	655	194	181	4		

- Molecule 15 is a protein called 50S ribosomal protein L15e.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	L15E	169	Total	C	N	O	S	0	0
			1423	899	283	236	5		

- Molecule 16 is a protein called 50S ribosomal protein L18e.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	L18E	112	Total	C	N	O	S	0	0
			895	576	163	153	3		

- Molecule 17 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	L18P	193	Total	C	N	O	S	0	0
			1539	990	274	274	1		

- Molecule 18 is a protein called 50S ribosomal protein L19e.

Mol	Chain	Residues	Atoms				AltConf	Trace
18	L19E	144	Total	C	N	O	0	0
			1206	753	247	206		

- Molecule 19 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	L22P	150	Total	C	N	O	S	0	0
			1223	782	225	213	3		

- Molecule 20 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	L23P	81	Total	C	N	O	S	0	0
			650	419	109	121	1		

- Molecule 21 is a protein called 50S ribosomal protein L24e.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	L24E	54	Total	C	N	O	S	0	0
			441	282	80	73	6		

- Molecule 22 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	L24P	122	Total	C	N	O	S	0	0
			989	620	189	176	4		

- Molecule 23 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	L29P	63	Total	C	N	O	S	0	0
			513	319	95	96	3		

- Molecule 24 is a protein called 50S ribosomal protein L30e.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	L30E	94	Total	C	N	O	S	0	0
			729	474	116	136	3		

- Molecule 25 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	L30P	155	Total	C	N	O	S	0	0
			1254	804	222	223	5		

- Molecule 26 is a protein called 50S ribosomal protein L31e.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	L31E	75	Total	C	N	O	S	0	0
			625	398	126	97	4		

- Molecule 27 is a protein called 50S ribosomal protein L32e.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	L32E	123	Total	C	N	O	S	0	0
			1010	650	193	166	1		

- Molecule 28 is a protein called 50S ribosomal protein L34e.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	L34E	77	Total	C	N	O	S	0	0
			629	395	119	110	5		

- Molecule 29 is a protein called 50S ribosomal protein L37Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	L37A	65	Total	C	N	O	S	0	0
			527	335	99	87	6		

- Molecule 30 is a protein called 50S ribosomal protein L37e.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	L37E	54	Total	C	N	O	S	0	0
			436	267	94	69	6		

- Molecule 31 is a protein called 50S ribosomal protein L39e.

Mol	Chain	Residues	Atoms				AltConf	Trace
31	L39E	49	Total	C	N	O	0	0
			414	265	88	61		

- Molecule 32 is a protein called 50S ribosomal protein L40E.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	L40E	55	Total	C	N	O	S	0	0
			439	273	89	72	5		

- Molecule 33 is a protein called 50S ribosomal protein L44e.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	L44E	92	Total	C	N	O	S	0	0
			753	474	144	129	6		

- Molecule 34 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	L7A1	123	Total	C	N	O	S	0	0
			935	593	155	184	3		
34	L7A2	123	Total	C	N	O	S	0	0
			935	593	155	184	3		
34	SL7A	123	Total	C	N	O	S	0	0
			935	593	155	184	3		

- Molecule 35 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	L15P	94	Total	C	N	O	S	0	0
			752	487	131	133	1		

- Molecule 36 is a protein called 50S ribosomal protein L21e.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	L21E	97	Total	C	N	O	S	0	0
			785	502	152	129	2		

- Molecule 37 is a protein called DUF2280 domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	L45A	101	Total	C	N	O	S	0	0
			816	515	141	156	4		

- Molecule 38 is a protein called Conserved protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	L46A	70	Total	C	N	O	S	0	0
			586	382	101	102	1		

- Molecule 39 is a protein called 50S ribosomal protein L47A.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	L47A	80	Total	C	N	O	S	0	0
			648	405	113	128	2		

- Molecule 40 is a protein called 50S ribosomal protein L1.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	AL1P	216	Total	C	N	O	S	0	0
			1715	1096	303	312	4		

- Molecule 41 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	AS3P	201	Total	C	N	O	S	0	0
			1576	1020	274	278	4		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
42	AS7P	193	Total	C	N	O	S	0	0
			1537	969	285	279	4		

- Molecule 43 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	S10P	100	Total	C	N	O	S	0	0
			824	522	154	142	6		

- Molecule 44 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	S13P	147	Total	C	N	O	S	0	0
			1204	753	230	217	4		

- Molecule 45 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	S14P	52	Total	C	N	O	S	0	0
			432	273	85	69	5		

- Molecule 46 is a protein called 30S ribosomal protein S17e.

Mol	Chain	Residues	Atoms				AltConf	Trace
46	S17E	62	Total	C	N	O	0	0
			517	326	92	99		

- Molecule 47 is a protein called 30S ribosomal protein S19e.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	S19E	149	Total	C	N	O	S	0	0
			1230	795	221	212	2		

- Molecule 48 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	S19P	115	Total	C	N	O	S	0	0
			968	620	181	162	5		

- Molecule 49 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	AS9P	136	Total	C	N	O	S	0	0
			1096	692	200	197	7		

- Molecule 50 is a protein called 30S ribosomal protein S28e.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	S28E	63	Total	C	N	O		0	0
			498	308	99	91			

- Molecule 51 is a protein called 30S ribosomal protein S27ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	S27A	54	Total	C	N	O	S	0	0
			435	274	79	76	6		

- Molecule 52 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	AS2P	196	Total	C	N	O	S	0	0
			1587	1022	277	286	2		

- Molecule 53 is a protein called 30S ribosomal protein S4e.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	AS4E	240	Total	C	N	O	S	0	0
			1925	1238	335	348	4		

- Molecule 54 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	AS4P	166	Total	C	N	O	S	0	0
			1370	874	252	241	3		

- Molecule 55 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	AS5P	204	Total	C	N	O	S	0	0
			1600	1028	277	287	8		

- Molecule 56 is a protein called 30S ribosomal protein S6e.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	AS6E	105	Total	C	N	O	S	0	0
			805	506	149	147	3		

- Molecule 57 is a protein called 30S ribosomal protein S8e.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	AS8E	126	Total	C	N	O	S	0	0
			993	619	187	187			

- Molecule 58 is a protein called Small ribosomal subunit protein uS8.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	AS8P	130	Total	C	N	O	S	0	0
			1028	661	181	182	4		

- Molecule 59 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	S11P	128	Total	C	N	O	S	0	0
			960	595	190	173	2		

- Molecule 60 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	S12P	143	Total	C	N	O	S	0	0
			1103	701	209	189	4		

- Molecule 61 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	S15P	149	Total	C	N	O	S	0	0
			1225	778	228	214	5		

- Molecule 62 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	S17P	111	Total	C	N	O	S	0	0
			885	557	165	160	3		

- Molecule 63 is a protein called 30S ribosomal protein S24e.

Mol	Chain	Residues	Atoms				AltConf	Trace
63	S24E	96	Total	C	N	O	0	0
			759	479	133	147		

- Molecule 64 is a protein called 30S ribosomal protein S27e.

Mol	Chain	Residues	Atoms					AltConf	Trace
64	S27E	59	Total	C	N	O	S	0	0
			458	294	83	76	5		

- Molecule 65 is a protein called 30S ribosomal protein S3Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
65	S3AE	189	Total	C	N	O	S	0	0
			1545	1004	264	276	1		

- Molecule 66 is a RNA chain called tRNA (76-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
66	APTN	76	Total	C	N	O	P	0	0
			1619	723	290	531	75		
66	AETN	76	Total	C	N	O	P	0	0
			1619	723	290	531	75		

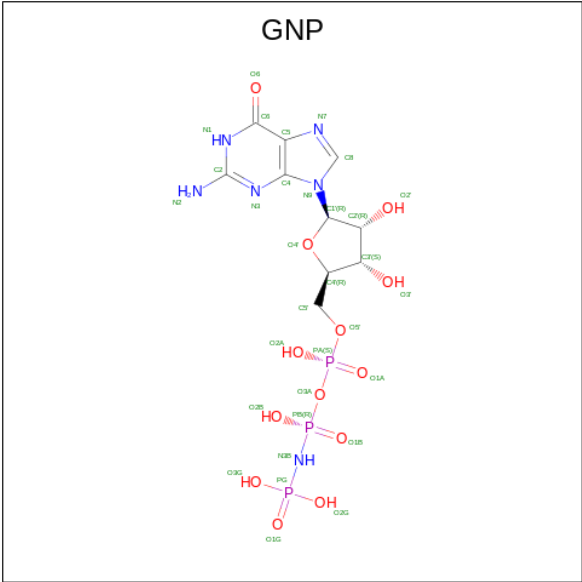
- Molecule 67 is a RNA chain called mRNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
67	AMRN	13	Total	C	N	O	P	0	0
			260	117	26	104	13		

- Molecule 68 is a protein called PHE-PHE-PHE-PHE-PHE-PHE.

Mol	Chain	Residues	Atoms				AltConf	Trace
68	APTP	6	Total	C	N	O	0	0
			67	54	6	7		

- Molecule 69 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).



Mol	Chain	Residues	Atoms					AltConf
69	AEFG	1	Total	C	N	O	P	0
			32	10	6	13	3	

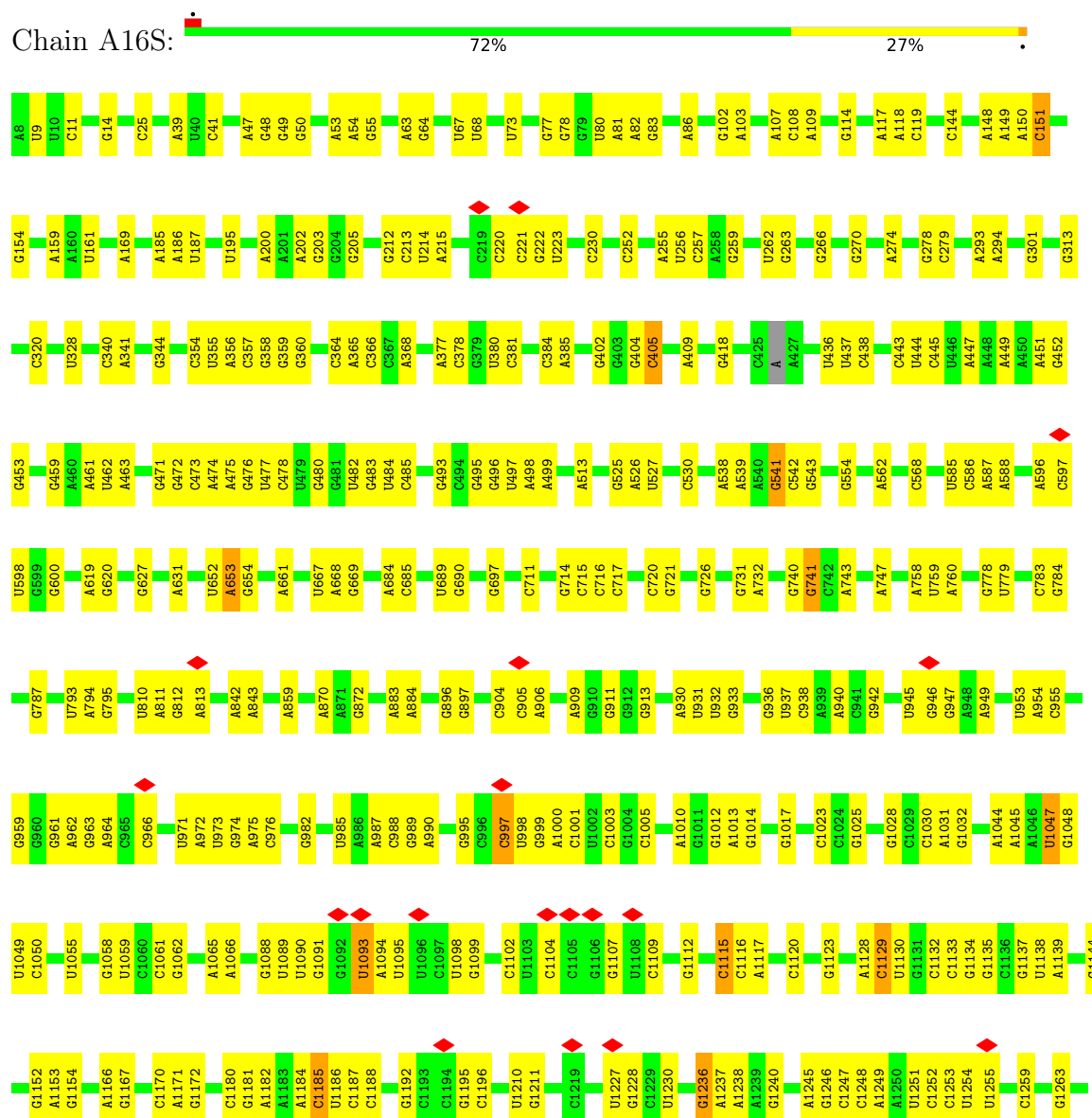
- Molecule 70 is UNKNOWN LIGAND (three-letter code: UNL) (formula:) (labeled as "Ligand of Interest" by depositor).

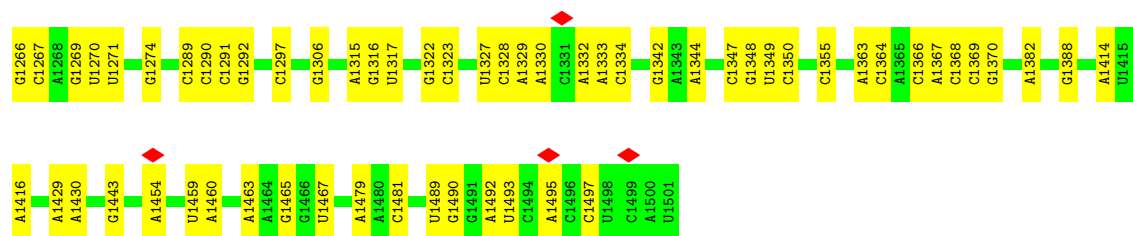
Mol	Chain	Residues	Atoms				AltConf
70	AS2P	34	Total	C	N	O	0
			171	102	34	35	
70	AS5P	16	Total	C	N	O	0
			80	48	16	16	
70	AS8P	7	Total	C	N	O	0
			35	21	7	7	

3 Residue-property plots

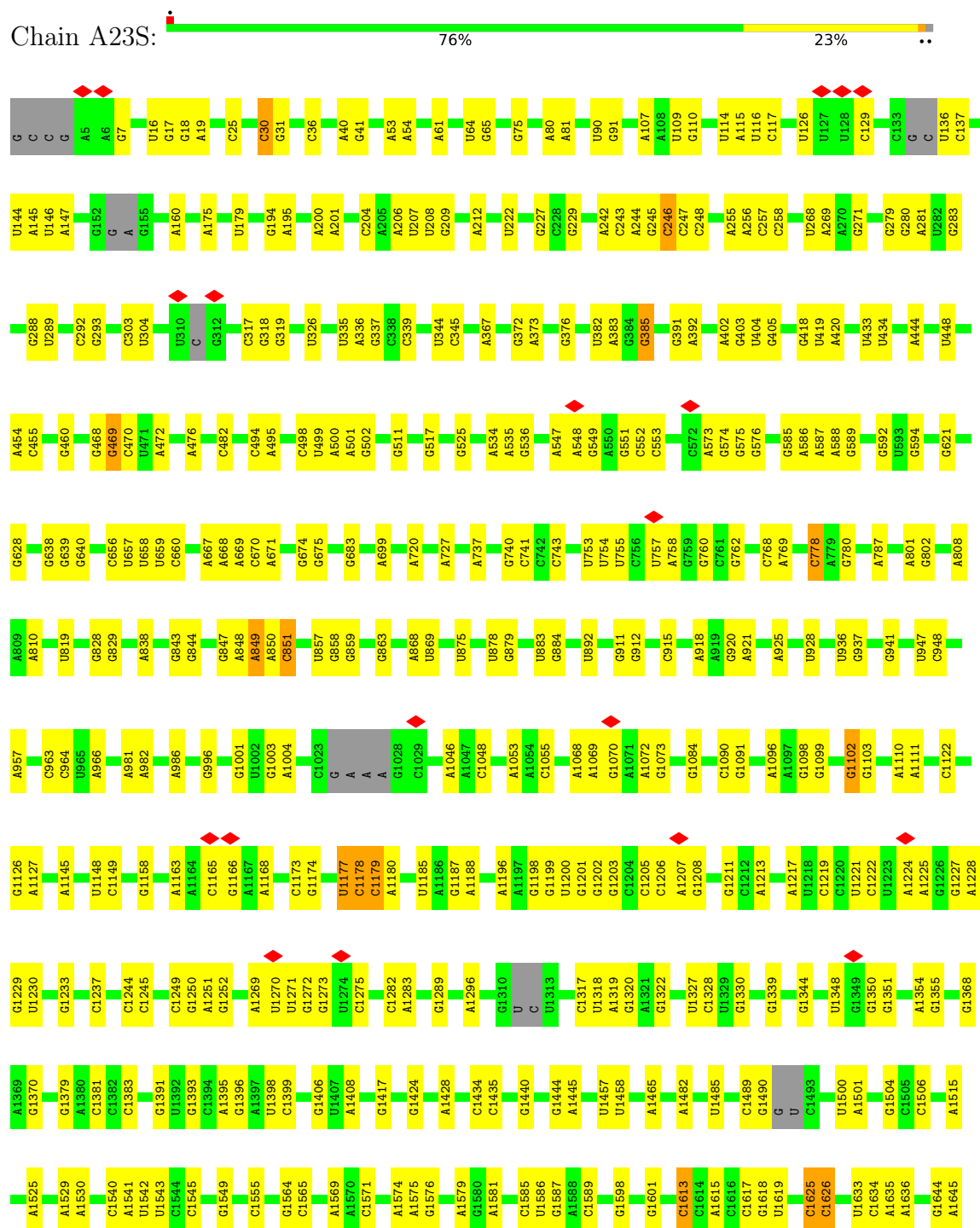
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: 16s rRNA (1493-MER)



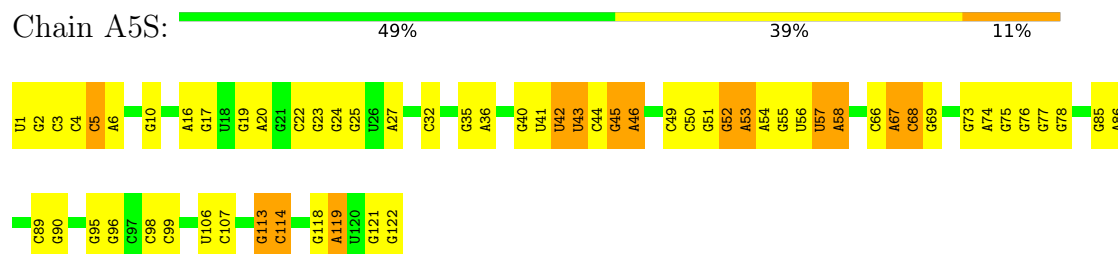


• Molecule 2: 23s rRNA (3000-MER)

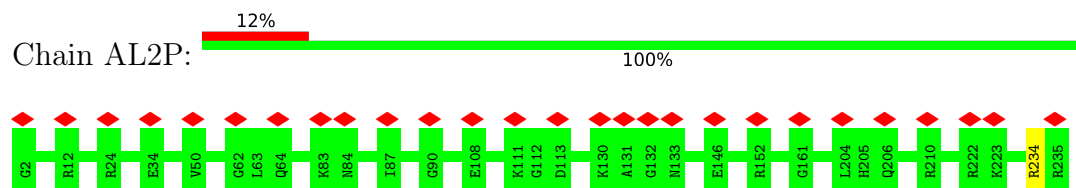




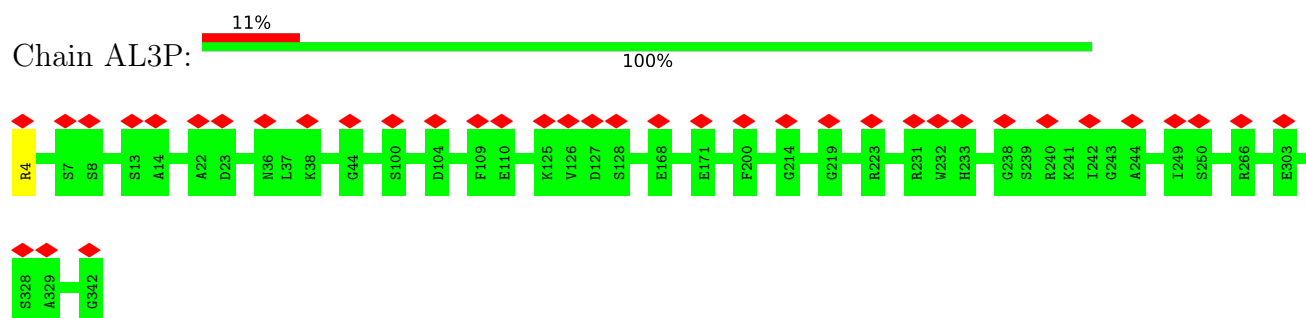
- Molecule 4: 5s rRNA (122-MER)



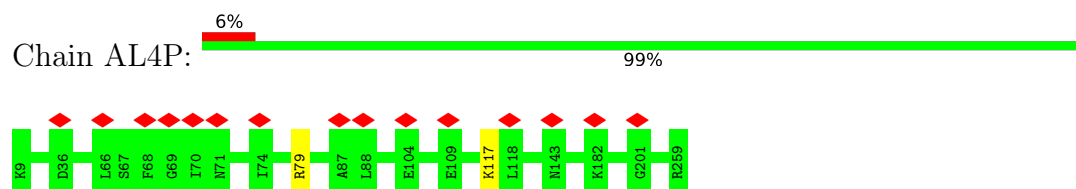
- Molecule 5: 50S ribosomal protein L2



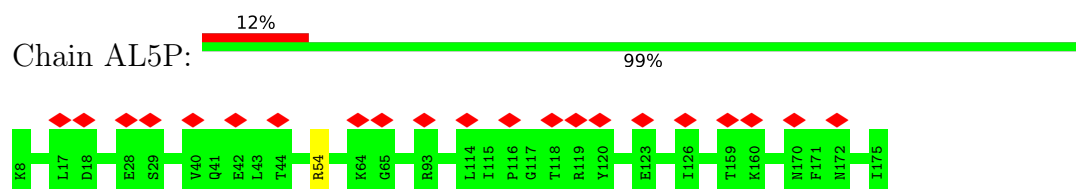
- Molecule 6: 50S ribosomal protein L3



- Molecule 7: 50S ribosomal protein L4



- Molecule 8: 50S ribosomal protein L5

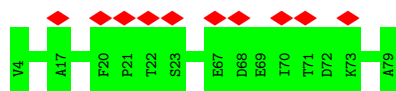


- Molecule 9: 50S ribosomal protein L6

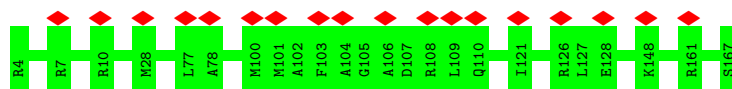




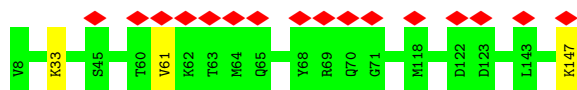
- Molecule 10: 50S ribosomal protein L18Ae



- Molecule 11: 50S ribosomal protein L10e



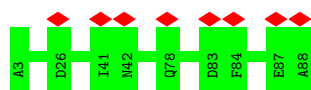
- Molecule 12: 50S ribosomal protein L13



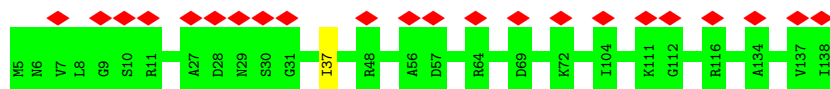
- Molecule 13: 50S ribosomal protein L14e



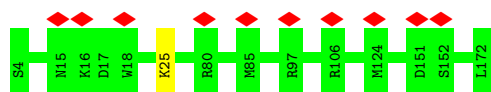
- Molecule 13: 50S ribosomal protein L14e



- Molecule 14: 50S ribosomal protein L14



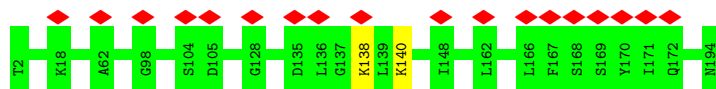
- Molecule 15: 50S ribosomal protein L15e



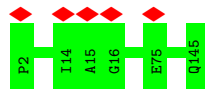
- Molecule 16: 50S ribosomal protein L18e



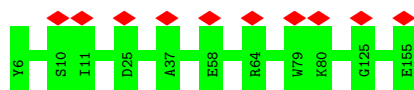
- Molecule 17: 50S ribosomal protein L18



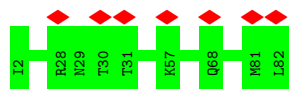
- Molecule 18: 50S ribosomal protein L19e



- Molecule 19: 50S ribosomal protein L22



- Molecule 20: 50S ribosomal protein L23

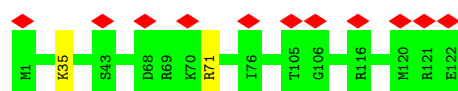


- Molecule 21: 50S ribosomal protein L24e

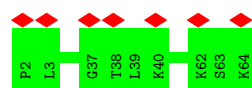


There are no outlier residues recorded for this chain.

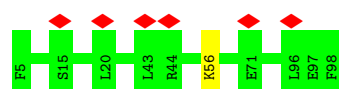
- Molecule 22: 50S ribosomal protein L24



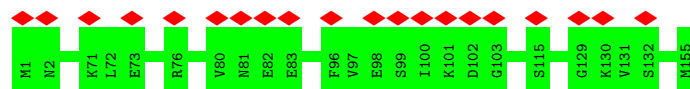
- Molecule 23: 50S ribosomal protein L29



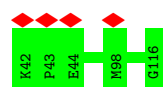
- Molecule 24: 50S ribosomal protein L30e



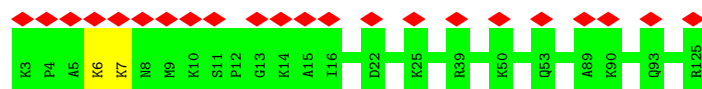
- Molecule 25: 50S ribosomal protein L30



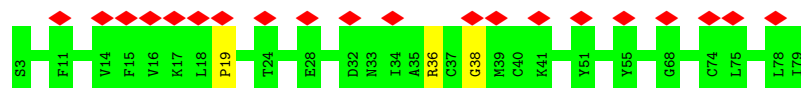
- Molecule 26: 50S ribosomal protein L31e



- Molecule 27: 50S ribosomal protein L32e

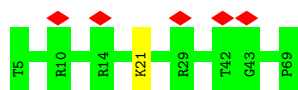


- Molecule 28: 50S ribosomal protein L34e



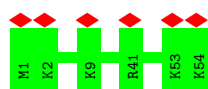
- Molecule 29: 50S ribosomal protein L37Ae

Chain L37A:  8% 98%



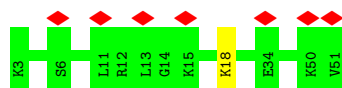
- Molecule 30: 50S ribosomal protein L37e

Chain L37E:  11% 100%

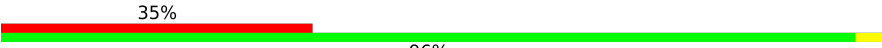


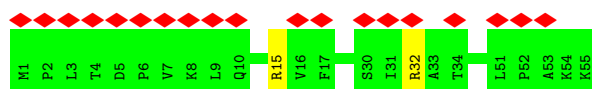
- Molecule 31: 50S ribosomal protein L39e

Chain L39E:  14% 98%



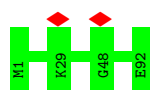
- Molecule 32: 50S ribosomal protein L40E

Chain L40E:  35% 96%



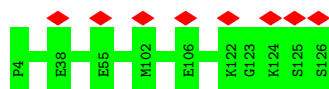
- Molecule 33: 50S ribosomal protein L44e

Chain L44E:  0% 100%



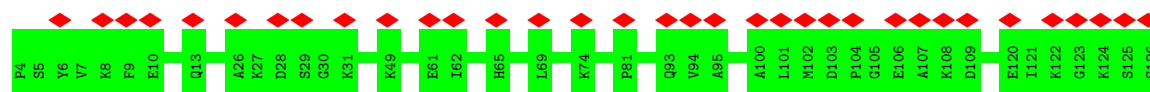
- Molecule 34: 50S ribosomal protein L7Ae

Chain L7A1:  7% 100%

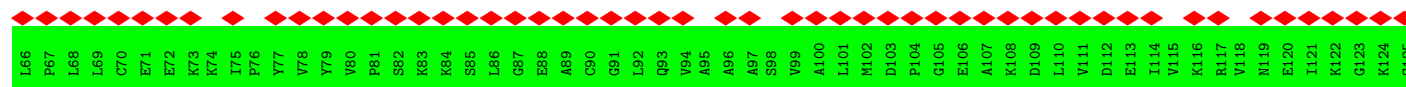
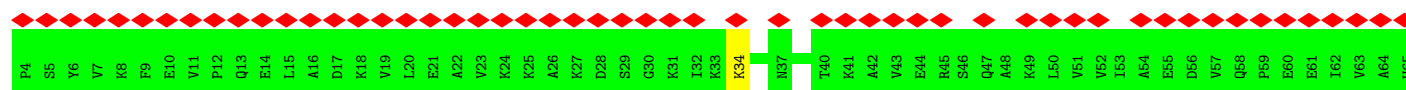


- Molecule 34: 50S ribosomal protein L7Ae

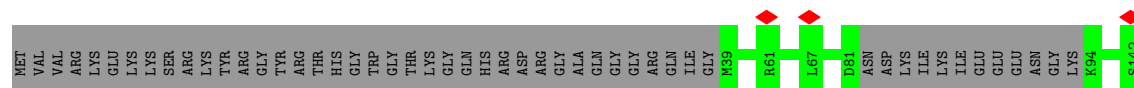
Chain L7A2:  28% 100%



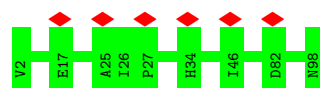
- Molecule 34: 50S ribosomal protein L7Ae



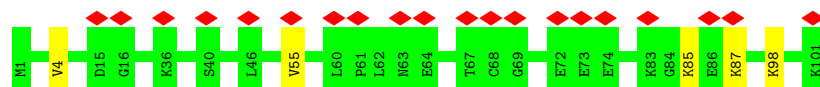
- Molecule 35: 50S ribosomal protein L15



- Molecule 36: 50S ribosomal protein L21e

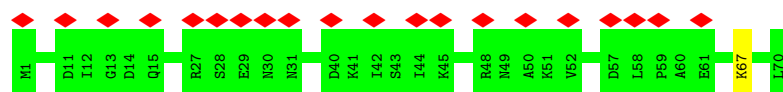


- Molecule 37: DUF2280 domain-containing protein



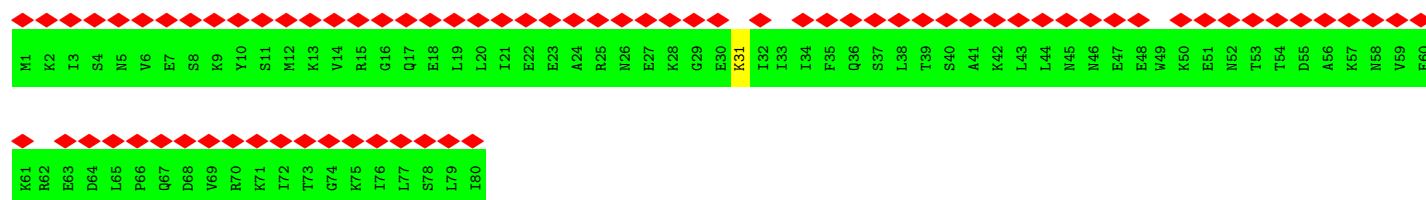
- Molecule 38: Conserved protein





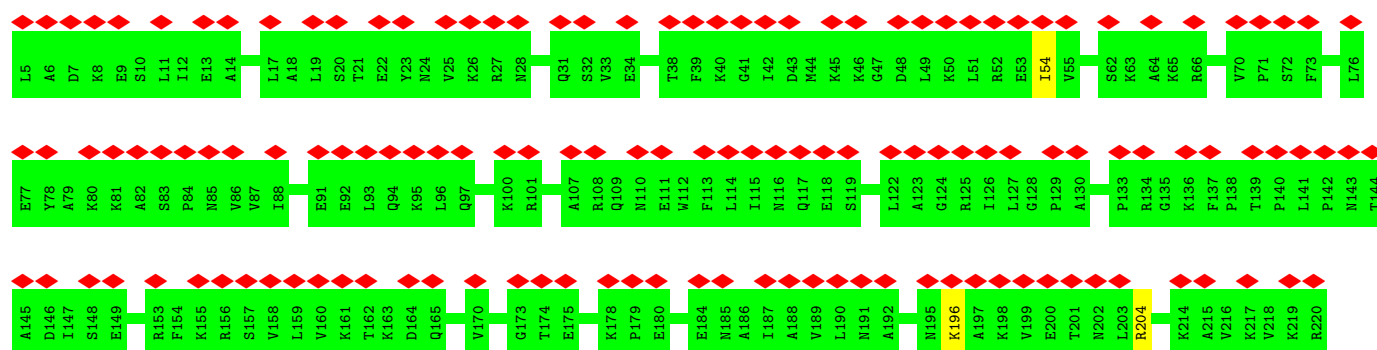
- Molecule 39: 50S ribosomal protein L47A

Chain L47A: 95%
99%



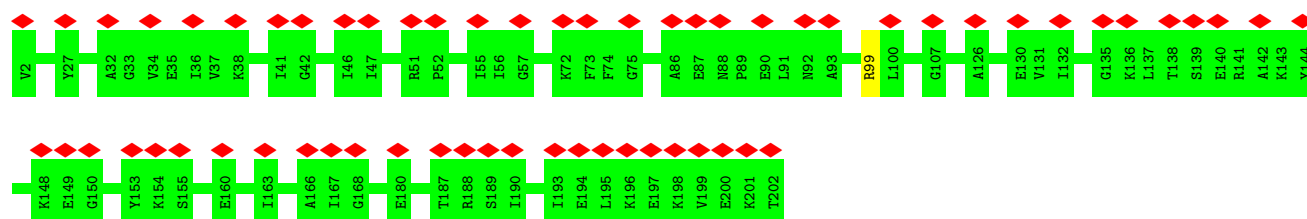
- Molecule 40: 50S ribosomal protein L1

Chain AL1P: 63%
99%



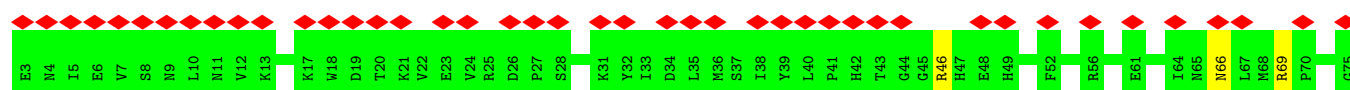
- Molecule 41: 30S ribosomal protein S3

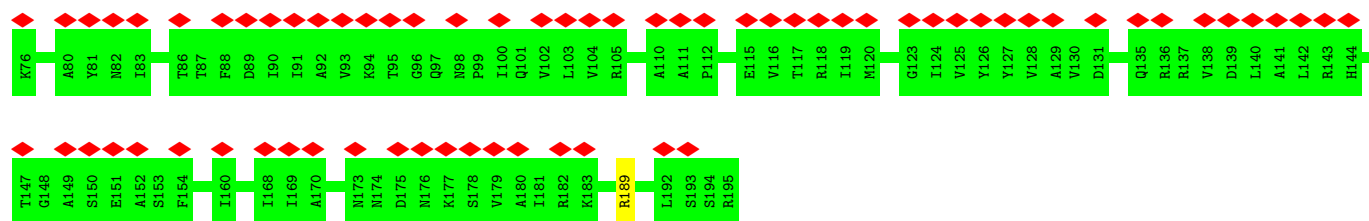
Chain AS3P: 30%
100%



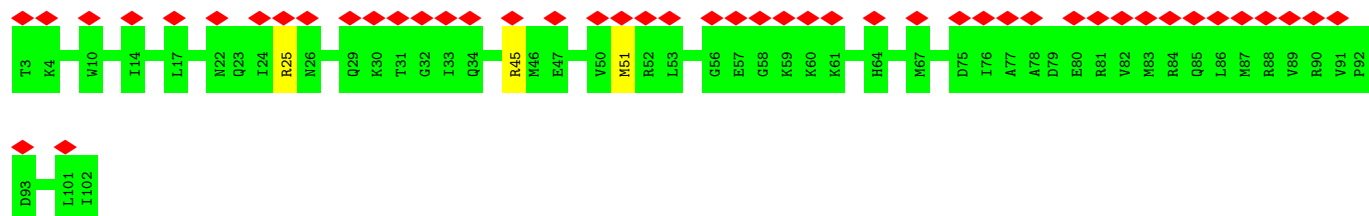
- Molecule 42: 30S ribosomal protein S7

Chain AS7P: 58%
98%

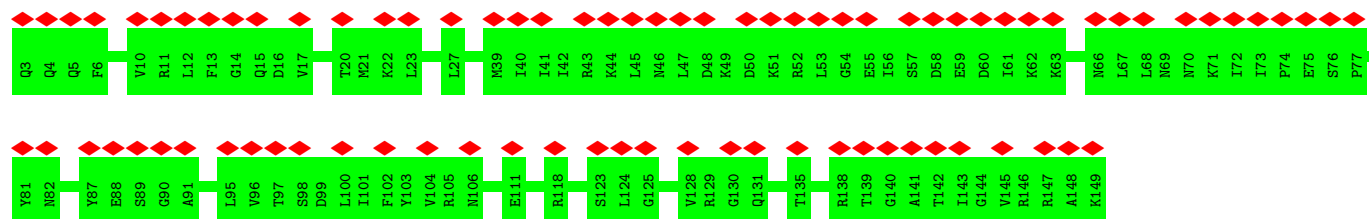




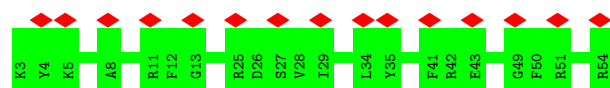
- Molecule 43: 30S ribosomal protein S10



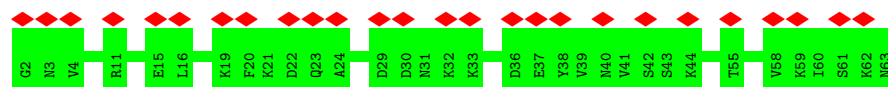
- Molecule 44: 30S ribosomal protein S13



- Molecule 45: 30S ribosomal protein S14 type Z

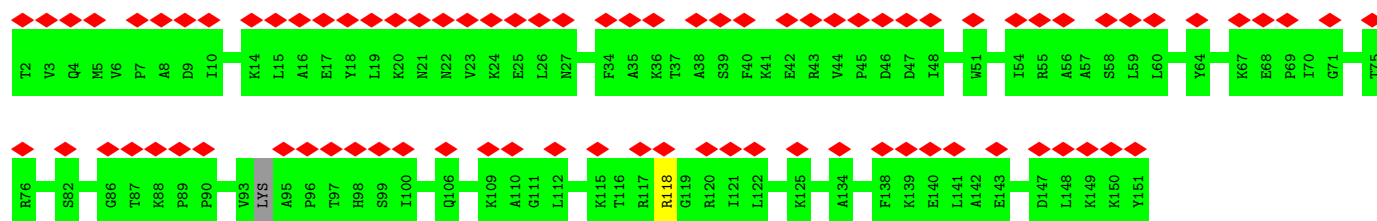


- Molecule 46: 30S ribosomal protein S17e



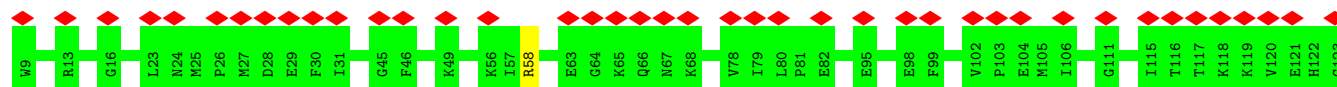
- Molecule 47: 30S ribosomal protein S19e





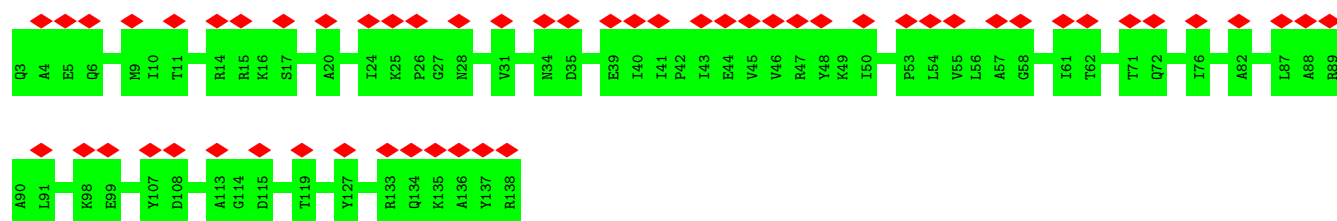
- Molecule 48: 30S ribosomal protein S19

Chain S19P: 36% 99%



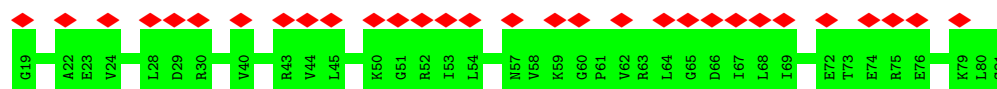
- Molecule 49: 30S ribosomal protein S9

Chain AS9P: 40% 100%



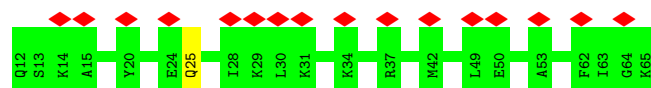
- Molecule 50: 30S ribosomal protein S28e

Chain S28E: 48% 100%



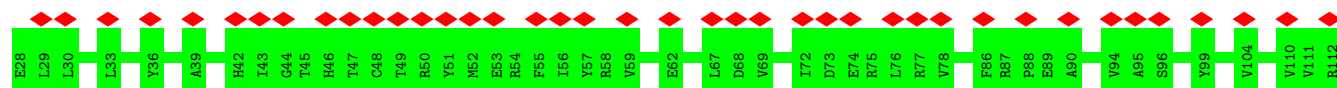
- Molecule 51: 30S ribosomal protein S27ae

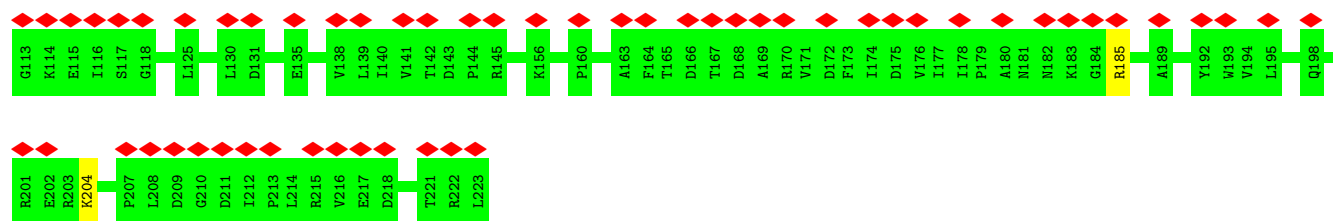
Chain S27A: 30% 98%



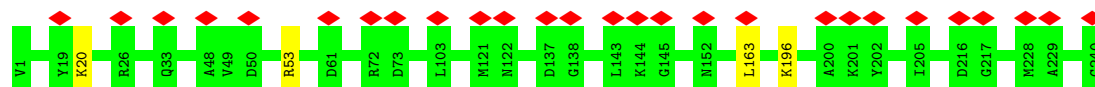
- Molecule 52: 30S ribosomal protein S2

Chain AS2P: 49% 99%

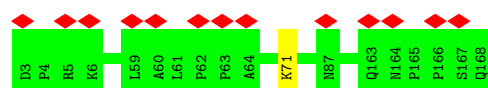




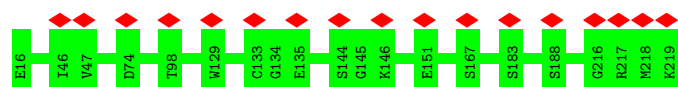
- Molecule 53: 30S ribosomal protein S4e



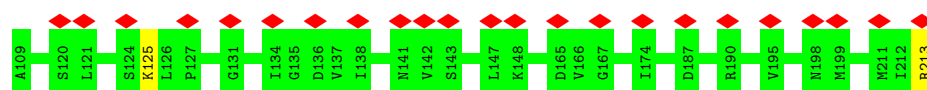
- Molecule 54: 30S ribosomal protein S4



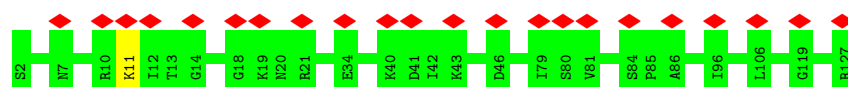
- Molecule 55: 30S ribosomal protein S5



- Molecule 56: 30S ribosomal protein S6e

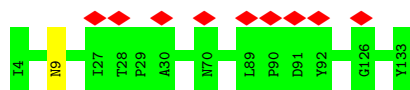


- Molecule 57: 30S ribosomal protein S8e

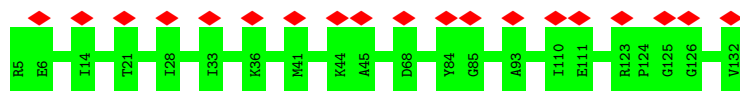


- Molecule 58: Small ribosomal subunit protein uS8





- Molecule 59: 30S ribosomal protein S11



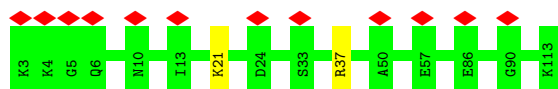
- Molecule 60: 30S ribosomal protein S12



- Molecule 61: 30S ribosomal protein S15



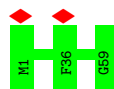
- Molecule 62: 30S ribosomal protein S17



- Molecule 63: 30S ribosomal protein S24e

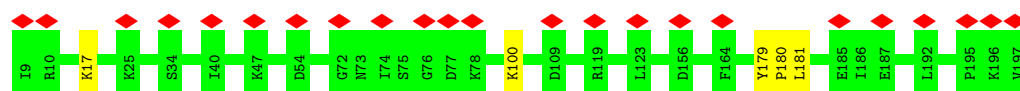


- Molecule 64: 30S ribosomal protein S27e




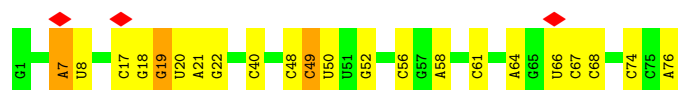
- Molecule 65: 30S ribosomal protein S3Ae

Chain S3AE:  12% 97% .



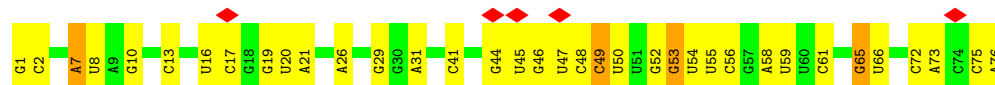
- Molecule 66: tRNA (76-MER)

Chain APTN:  71% 25% .



- Molecule 66: tRNA (76-MER)

Chain AETN:  7% 53% 42% 5%




- Molecule 67: mRNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

Chain AMRN:  23% 46% 54%



- Molecule 68: PHE-PHE-PHE-PHE-PHE-PHE

Chain APTP:  83% 100%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	3841	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 ($e^-/\text{\AA}^2$)	26.7	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	1.282	Depositor
Minimum map value	-0.504	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.100	Depositor
Recommended contour level	0.33	Depositor
Map size (Å)	413.06, 413.06, 413.06	wwPDB
Map dimensions	380, 380, 380	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.087, 1.087, 1.087	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	A16S	0.56	6/35889 (0.0%)	0.90	61/56002 (0.1%)
2	A23S	0.43	1/72029 (0.0%)	0.88	109/112378 (0.1%)
3	AEFG	0.28	0/5771	0.54	1/7808 (0.0%)
4	A5S	0.34	0/2917	0.70	0/4549
5	AL2P	0.30	0/1787	0.55	0/2409
6	AL3P	0.29	0/2758	0.53	0/3727
7	AL4P	0.29	0/1956	0.55	0/2635
8	AL5P	0.28	0/1364	0.57	0/1827
9	AL6P	0.29	0/1450	0.49	0/1949
10	ALX0	0.32	0/638	0.62	0/851
11	L10E	0.30	0/1334	0.55	0/1787
12	L13P	0.28	0/1123	0.58	0/1502
13	L141	0.27	0/673	0.51	0/900
13	L142	0.27	0/673	0.51	0/900
14	L14P	0.30	0/1054	0.56	0/1425
15	L15E	0.32	0/1458	0.58	0/1956
16	L18E	0.28	0/907	0.51	0/1214
17	L18P	0.29	0/1570	0.49	0/2115
18	L19E	0.27	0/1223	0.54	0/1622
19	L22P	0.29	0/1246	0.50	0/1671
20	L23P	0.29	0/655	0.51	0/874
21	L24E	0.29	0/451	0.50	0/599
22	L24P	0.30	0/1000	0.57	0/1329
23	L29P	0.25	0/513	0.54	0/678
24	L30E	0.27	0/738	0.49	0/985
25	L30P	0.30	0/1278	0.55	0/1713
26	L31E	0.28	0/632	0.58	0/837
27	L32E	0.28	0/1027	0.53	0/1366
28	L34E	0.31	0/642	0.70	0/854
29	L37A	0.31	0/542	0.61	0/726
30	L37E	0.31	0/445	0.57	0/585
31	L39E	0.26	0/422	0.60	0/562

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	L40E	0.27	0/443	0.63	0/587
33	L44E	0.28	0/763	0.55	0/1008
34	L7A1	0.27	0/946	0.43	0/1272
34	L7A2	0.25	0/946	0.43	0/1272
34	SL7A	0.25	0/946	0.45	0/1272
35	L15P	0.28	0/766	0.49	0/1023
36	L21E	0.29	0/800	0.51	0/1067
37	L45A	0.27	0/824	0.56	0/1094
38	L46A	0.26	0/595	0.47	0/793
39	L47A	0.24	0/652	0.52	0/870
40	AL1P	0.26	0/1739	0.52	0/2338
41	AS3P	0.27	0/1599	0.51	0/2147
42	AS7P	0.26	0/1561	0.55	0/2105
43	S10P	0.25	0/840	0.55	0/1132
44	S13P	0.26	0/1221	0.55	0/1634
45	S14P	0.27	0/441	0.61	0/583
46	S17E	0.25	0/523	0.47	0/696
47	S19E	0.25	0/1257	0.51	0/1691
48	S19P	0.25	0/985	0.56	0/1310
49	AS9P	0.26	0/1115	0.54	0/1496
50	S28E	0.24	0/500	0.60	0/669
51	S27A	0.30	0/444	0.62	0/590
52	AS2P	0.27	0/1621	0.55	0/2202
53	AS4E	0.28	0/1956	0.55	0/2635
54	AS4P	0.28	0/1399	0.55	0/1883
55	AS5P	0.28	0/1631	0.52	0/2200
56	AS6E	0.28	0/815	0.57	0/1093
57	AS8E	0.27	0/1005	0.56	0/1342
58	AS8P	0.30	0/1046	0.57	0/1410
59	S11P	0.27	0/976	0.59	0/1315
60	S12P	0.29	0/1120	0.58	0/1495
61	S15P	0.28	0/1250	0.54	0/1677
62	S17P	0.28	0/899	0.56	0/1203
63	S24E	0.26	0/769	0.50	0/1034
64	S27E	0.27	0/465	0.49	0/618
65	S3AE	0.28	0/1573	0.54	0/2115
66	AETN	0.53	2/1809 (0.1%)	2.13	25/2819 (0.9%)
66	APTN	0.51	3/1809 (0.2%)	1.52	25/2819 (0.9%)
67	AMRN	0.25	0/285	0.86	0/438
68	APTP	0.33	0/72	0.34	0/93
All	All	0.41	12/186571 (0.0%)	0.81	221/275375 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	AEFG	0	1
12	L13P	0	1
28	L34E	0	2
42	AS7P	0	1
51	S27A	0	1
53	AS4E	0	1
61	S15P	0	1
65	S3AE	0	2
All	All	0	10

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A16S	1236	G	C6-N1	37.36	1.65	1.39
1	A16S	1236	G	N3-C4	34.54	1.59	1.35
1	A16S	1236	G	N1-C2	33.69	1.64	1.37
1	A16S	1236	G	C2-N3	32.36	1.58	1.32
1	A16S	1236	G	C5-C4	28.52	1.58	1.38
1	A16S	1236	G	C5-C6	22.22	1.64	1.42
66	AETN	61	C	C2-N3	-7.07	1.30	1.35
66	APTN	7	A	N9-C4	-6.79	1.33	1.37
2	A23S	1179	C	N1-C6	-5.91	1.33	1.37
66	AETN	61	C	N3-C4	-5.39	1.30	1.33
66	APTN	19	G	N1-C2	-5.38	1.33	1.37
66	APTN	19	G	C6-N1	-5.04	1.36	1.39

All (221) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
66	AETN	61	C	N3-C2-O2	-49.60	87.18	121.90
66	AETN	61	C	N1-C2-O2	47.86	147.62	118.90
66	APTN	19	G	N3-C2-N2	33.59	143.42	119.90
66	APTN	19	G	N1-C2-N2	-32.19	87.23	116.20
66	AETN	29	G	N1-C2-N2	-28.30	90.73	116.20
66	AETN	65	G	N1-C2-N2	-27.17	91.75	116.20
66	AETN	29	G	N3-C2-N2	25.06	137.44	119.90
66	AETN	65	G	N3-C2-N2	21.28	134.80	119.90
2	A23S	1586	U	C5-C6-N1	18.71	132.05	122.70
66	AETN	61	C	C6-N1-C2	-17.09	113.47	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A16S	1236	G	C2-N3-C4	16.61	120.20	111.90
66	APTN	19	G	C5-C6-O6	16.01	138.21	128.60
66	AETN	65	G	N1-C2-N3	15.93	133.46	123.90
66	APTN	56	C	N1-C2-O2	15.65	128.29	118.90
66	AETN	61	C	C5-C4-N4	15.21	130.85	120.20
2	A23S	1179	C	C6-N1-C2	-13.44	114.92	120.30
66	AETN	29	G	N1-C2-N3	13.21	131.83	123.90
66	AETN	61	C	N3-C4-N4	-13.00	108.90	118.00
2	A23S	1179	C	C5-C6-N1	12.94	127.47	121.00
66	AETN	41	C	N1-C2-O2	12.50	126.40	118.90
66	APTN	19	G	N1-C6-O6	-12.43	112.44	119.90
66	APTN	19	G	C2-N3-C4	-12.31	105.75	111.90
66	AETN	49	C	N3-C2-O2	-12.21	113.35	121.90
66	APTN	56	C	N3-C2-O2	-11.81	113.63	121.90
66	AETN	41	C	N3-C2-O2	-11.72	113.69	121.90
66	AETN	65	G	C2-N3-C4	-11.71	106.04	111.90
1	A16S	1236	G	N3-C4-N9	11.62	132.97	126.00
1	A16S	1236	G	N1-C2-N3	-11.55	116.97	123.90
2	A23S	1586	U	C6-N1-C2	-11.39	114.17	121.00
2	A23S	469	G	N3-C4-N9	11.18	132.71	126.00
2	A23S	255	A	OP2-P-O3'	-11.17	80.63	105.20
2	A23S	469	G	C6-C5-N7	-11.00	123.80	130.40
66	APTN	7	A	N9-C4-C5	10.67	110.07	105.80
2	A23S	469	G	C8-N9-C1'	-10.62	113.19	127.00
2	A23S	469	G	N9-C4-C5	-10.60	101.16	105.40
2	A23S	246	C	N1-C2-O2	10.58	125.25	118.90
66	APTN	7	A	N3-C4-N9	-10.57	118.94	127.40
66	AETN	29	G	C2-N3-C4	-10.46	106.67	111.90
2	A23S	1585	C	C6-N1-C2	-10.40	116.14	120.30
2	A23S	469	G	C4-N9-C1'	10.27	139.85	126.50
2	A23S	469	G	C4-C5-N7	10.08	114.83	110.80
2	A23S	1179	C	N1-C2-O2	10.03	124.92	118.90
66	AETN	53	G	N3-C2-N2	9.63	126.64	119.90
66	AETN	49	C	N1-C2-O2	9.53	124.62	118.90
2	A23S	246	C	C2-N1-C1'	9.48	129.23	118.80
66	APTN	7	A	N1-C6-N6	-9.33	113.00	118.60
66	APTN	7	A	C5-C6-N6	9.26	131.11	123.70
2	A23S	246	C	C6-N1-C1'	-9.19	109.77	120.80
1	A16S	1236	G	N3-C4-C5	-9.19	124.01	128.60
66	APTN	19	G	N1-C2-N3	9.08	129.35	123.90
2	A23S	1586	U	C2-N1-C1'	9.00	128.50	117.70
2	A23S	255	A	OP1-P-O3'	-8.97	85.47	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A23S	1178	C	C2-N1-C1'	8.93	128.62	118.80
1	A16S	1236	G	C4-C5-N7	-8.78	107.29	110.80
1	A16S	1185	C	N1-C2-O2	8.61	124.07	118.90
66	AETN	61	C	N1-C2-N3	8.58	125.20	119.20
66	APTN	7	A	C6-C5-N7	8.40	138.18	132.30
2	A23S	1683	G	N3-C4-N9	-8.28	121.03	126.00
1	A16S	1047	U	N3-C2-O2	-8.24	116.43	122.20
66	APTN	7	A	N1-C2-N3	8.17	133.38	129.30
1	A16S	25	C	N3-C2-O2	-8.16	116.19	121.90
1	A16S	1185	C	C2-N1-C1'	8.16	127.78	118.80
66	APTN	7	A	C2-N3-C4	-8.15	106.53	110.60
1	A16S	1236	G	N7-C8-N9	8.07	117.14	113.10
2	A23S	1179	C	N3-C2-O2	-8.04	116.27	121.90
1	A16S	1180	C	N3-C2-O2	-8.04	116.27	121.90
2	A23S	1793	C	N3-C2-O2	-7.95	116.34	121.90
2	A23S	1457	U	C2-N1-C1'	7.94	127.22	117.70
66	AETN	53	G	C6-N1-C2	7.88	129.83	125.10
66	AETN	65	G	C6-N1-C2	-7.85	120.39	125.10
2	A23S	1625	C	O4'-C1'-N1	7.76	114.41	108.20
2	A23S	2641	C	N3-C2-O2	-7.61	116.58	121.90
2	A23S	3017	C	N1-C2-O2	7.48	123.39	118.90
1	A16S	1236	G	C5-C6-N1	7.43	115.21	111.50
66	APTN	7	A	C4-C5-N7	-7.33	107.03	110.70
66	APTN	56	C	C2-N3-C4	7.30	123.55	119.90
2	A23S	1613	C	N1-C2-O2	7.29	123.28	118.90
2	A23S	1683	G	C4-N9-C1'	-7.24	117.09	126.50
2	A23S	2022	C	N3-C2-O2	-7.22	116.85	121.90
1	A16S	1047	U	C2-N1-C1'	7.20	126.34	117.70
2	A23S	3017	C	N3-C2-O2	-7.17	116.88	121.90
2	A23S	1683	G	C8-N9-C1'	7.16	136.31	127.00
1	A16S	1047	U	N1-C2-O2	7.13	127.79	122.80
66	APTN	56	C	N3-C4-C5	-7.09	119.06	121.90
2	A23S	656	C	N3-C2-O2	-7.08	116.94	121.90
2	A23S	1685	C	N3-C2-O2	-7.05	116.97	121.90
66	APTN	7	A	C8-N9-C1'	7.03	140.36	127.70
2	A23S	1435	C	N3-C2-O2	-7.00	117.00	121.90
2	A23S	256	A	OP1-P-OP2	6.99	130.09	119.60
1	A16S	144	C	N3-C2-O2	-6.92	117.05	121.90
66	APTN	49	C	C2-N1-C1'	6.85	126.33	118.80
2	A23S	1585	C	C2-N1-C1'	6.85	126.33	118.80
2	A23S	1918	C	N3-C2-O2	-6.82	117.12	121.90
1	A16S	1093	U	C5-C4-O4	-6.79	121.82	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A23S	1586	U	N3-C4-O4	6.79	124.15	119.40
2	A23S	1178	C	C6-N1-C2	-6.78	117.59	120.30
2	A23S	1626	C	N3-C2-O2	-6.78	117.16	121.90
1	A16S	1003	C	C2-N1-C1'	6.75	126.23	118.80
1	A16S	741	G	N3-C4-N9	-6.67	122.00	126.00
1	A16S	720	C	C2-N1-C1'	6.66	126.13	118.80
66	APTN	7	A	C4-N9-C1'	-6.65	114.32	126.30
66	AETN	53	G	C5-C6-N1	-6.59	108.21	111.50
2	A23S	2090	C	N3-C2-O2	-6.57	117.30	121.90
1	A16S	1133	C	N3-C2-O2	-6.57	117.30	121.90
1	A16S	1185	C	C6-N1-C1'	-6.51	112.99	120.80
1	A16S	1003	C	N1-C2-O2	6.51	122.80	118.90
1	A16S	717	C	N3-C2-O2	-6.43	117.40	121.90
1	A16S	1012	G	N1-C6-O6	-6.39	116.06	119.90
1	A16S	1187	C	N3-C2-O2	-6.39	117.42	121.90
2	A23S	469	G	N3-C2-N2	6.35	124.35	119.90
3	AEFG	205	LYS	CB-CG-CD	-6.27	95.30	111.60
2	A23S	246	C	O4'-C1'-N1	6.24	113.19	108.20
2	A23S	279	G	C5-C6-O6	6.23	132.34	128.60
2	A23S	1626	C	N1-C2-O2	6.22	122.63	118.90
1	A16S	1236	G	N3-C2-N2	6.20	124.24	119.90
1	A16S	1093	U	O4'-C1'-N1	6.19	113.15	108.20
2	A23S	1457	U	N1-C2-O2	6.19	127.13	122.80
2	A23S	1625	C	C6-N1-C2	-6.17	117.83	120.30
1	A16S	354	C	N3-C2-O2	-6.14	117.60	121.90
2	A23S	1179	C	O4'-C1'-N1	6.14	113.11	108.20
2	A23S	1435	C	C6-N1-C2	-6.14	117.84	120.30
1	A16S	1093	U	N3-C4-O4	6.09	123.66	119.40
1	A16S	711	C	N3-C2-O2	-6.08	117.64	121.90
2	A23S	1683	G	C6-C5-N7	6.07	134.04	130.40
1	A16S	568	C	N3-C2-O2	-6.05	117.67	121.90
66	AETN	29	G	C6-N1-C2	-6.00	121.50	125.10
1	A16S	1187	C	N1-C2-O2	6.00	122.50	118.90
2	A23S	1178	C	C5-C6-N1	6.00	124.00	121.00
1	A16S	1012	G	C5-C6-O6	5.99	132.19	128.60
66	APTN	7	A	N3-C4-C5	5.99	130.99	126.80
2	A23S	2265	C	N1-C2-O2	5.96	122.48	118.90
1	A16S	154	G	C5-C6-O6	5.96	132.18	128.60
1	A16S	1132	C	N1-C2-O2	5.96	122.47	118.90
1	A16S	1236	G	N9-C4-C5	-5.95	103.02	105.40
1	A16S	1129	C	C2-N1-C1'	5.95	125.34	118.80
66	APTN	56	C	C6-N1-C2	-5.95	117.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A23S	1625	C	C6-N1-C1'	5.93	127.92	120.80
2	A23S	1802	C	C2-N1-C1'	5.92	125.32	118.80
2	A23S	2147	C	N1-C2-O2	5.92	122.45	118.90
2	A23S	1613	C	N3-C2-O2	-5.92	117.75	121.90
1	A16S	405	C	C2-N1-C1'	5.92	125.31	118.80
2	A23S	1823	G	N1-C6-O6	-5.91	116.35	119.90
2	A23S	1178	C	C6-N1-C1'	-5.89	113.73	120.80
2	A23S	1179	C	C5'-C4'-O4'	5.89	116.17	109.10
2	A23S	258	C	N3-C2-O2	-5.89	117.78	121.90
1	A16S	14	G	C5-C6-O6	5.88	132.13	128.60
2	A23S	303	C	N3-C2-O2	-5.88	117.79	121.90
1	A16S	151	C	N3-C2-O2	-5.87	117.79	121.90
2	A23S	1457	U	N3-C2-O2	-5.87	118.09	122.20
2	A23S	2090	C	N1-C2-O2	5.86	122.42	118.90
2	A23S	1178	C	N1-C2-O2	5.85	122.41	118.90
2	A23S	129	C	N3-C2-O2	-5.80	117.84	121.90
2	A23S	1728	G	C5-C6-O6	-5.78	125.13	128.60
2	A23S	1683	G	O4'-C1'-N9	5.76	112.81	108.20
2	A23S	1625	C	N1-C2-N3	5.76	123.23	119.20
1	A16S	14	G	N1-C6-O6	-5.76	116.45	119.90
2	A23S	1683	G	N1-C6-O6	-5.76	116.45	119.90
2	A23S	469	G	N1-C6-O6	5.75	123.35	119.90
1	A16S	1144	G	N3-C4-N9	5.72	129.43	126.00
2	A23S	303	C	N1-C2-O2	5.72	122.33	118.90
2	A23S	1625	C	C5-C4-N4	5.71	124.20	120.20
1	A16S	716	C	C2-N1-C1'	5.65	125.02	118.80
66	AETN	1	G	N3-C4-N9	5.65	129.39	126.00
2	A23S	1613	C	C2-N1-C1'	5.64	125.00	118.80
2	A23S	1179	C	C2-N1-C1'	5.62	124.98	118.80
2	A23S	1090	C	N3-C2-O2	-5.60	117.98	121.90
2	A23S	1802	C	N1-C2-O2	5.60	122.26	118.90
2	A23S	1585	C	C5-C6-N1	5.58	123.79	121.00
2	A23S	469	G	C5-C6-O6	-5.57	125.26	128.60
2	A23S	2692	G	N1-C2-N2	-5.57	111.19	116.20
2	A23S	1103	G	N3-C4-N9	-5.54	122.67	126.00
2	A23S	1683	G	C5-C6-O6	5.54	131.92	128.60
2	A23S	482	C	N1-C2-O2	5.52	122.21	118.90
2	A23S	1683	G	N3-C4-C5	5.52	131.36	128.60
2	A23S	1177	U	C2-N1-C1'	5.52	124.32	117.70
2	A23S	1683	G	N9-C4-C5	5.51	107.60	105.40
1	A16S	154	G	N1-C6-O6	-5.51	116.59	119.90
2	A23S	778	C	N1-C2-O2	5.50	122.20	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
66	APT	56	C	C5-C6-N1	5.48	123.74	121.00
1	A16S	1180	C	N1-C2-N3	5.48	123.04	119.20
2	A23S	2225	C	C2-N1-C1'	5.47	124.82	118.80
2	A23S	2265	C	N3-C2-O2	-5.45	118.09	121.90
2	A23S	279	G	N1-C6-O6	-5.42	116.64	119.90
1	A16S	720	C	C6-N1-C1'	-5.42	114.30	120.80
2	A23S	30	C	C2-N1-C1'	5.41	124.75	118.80
1	A16S	653	A	P-O3'-C3'	5.38	126.16	119.70
2	A23S	1457	U	C6-N1-C1'	-5.36	113.69	121.20
2	A23S	279	G	N3-C4-N9	-5.36	122.79	126.00
2	A23S	2147	C	C2-N1-C1'	5.36	124.69	118.80
2	A23S	339	C	N3-C2-O2	-5.35	118.15	121.90
2	A23S	1586	U	C5-C4-O4	-5.34	122.70	125.90
2	A23S	2022	C	N1-C2-O2	5.33	122.09	118.90
66	APT	49	C	C6-N1-C1'	-5.28	114.46	120.80
1	A16S	717	C	C6-N1-C2	-5.27	118.19	120.30
1	A16S	955	C	N3-C2-O2	-5.26	118.22	121.90
2	A23S	1102	G	N3-C4-N9	5.25	129.15	126.00
2	A23S	2308	C	N3-C2-O2	-5.20	118.26	121.90
1	A16S	741	G	C4-N9-C1'	-5.19	119.75	126.50
2	A23S	849	A	O4'-C1'-N9	5.18	112.35	108.20
2	A23S	385	G	N1-C6-O6	-5.17	116.80	119.90
2	A23S	1434	C	N1-C2-O2	5.17	122.00	118.90
2	A23S	2119	C	N3-C2-O2	-5.16	118.28	121.90
2	A23S	851	C	N3-C2-O2	-5.16	118.29	121.90
1	A16S	716	C	C6-N1-C1'	-5.15	114.62	120.80
2	A23S	2692	G	C5-C6-O6	5.15	131.69	128.60
1	A16S	80	U	C2-N1-C1'	5.13	123.86	117.70
1	A16S	541	G	P-O3'-C3'	5.12	125.84	119.70
1	A16S	741	G	N3-C4-C5	5.11	131.16	128.60
2	A23S	1625	C	N3-C2-O2	-5.10	118.33	121.90
1	A16S	997	C	C2-N1-C1'	5.09	124.40	118.80
66	AET	7	A	C5-C6-N1	5.08	120.24	117.70
2	A23S	1585	C	N3-C2-O2	-5.07	118.35	121.90
2	A23S	248	C	C2-N1-C1'	5.05	124.35	118.80
2	A23S	1918	C	N1-C2-O2	5.04	121.93	118.90
1	A16S	1115	C	N1-C2-O2	5.04	121.92	118.90
1	A16S	1003	C	N3-C2-O2	-5.04	118.37	121.90
1	A16S	741	G	N3-C2-N2	-5.03	116.38	119.90
2	A23S	469	G	N3-C4-C5	-5.02	126.09	128.60
1	A16S	1180	C	C5-C4-N4	5.02	123.72	120.20
1	A16S	1115	C	N3-C2-O2	-5.01	118.39	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A16S	1023	C	C2-N1-C1'	5.01	124.31	118.80

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	AEFG	343	TYR	Peptide
53	AS4E	163	LEU	Peptide
42	AS7P	69	ARG	Peptide
12	L13P	61	VAL	Peptide
28	L34E	36	ARG	Peptide
28	L34E	38	GLY	Peptide
61	S15P	20	PRO	Peptide
51	S27A	25	GLN	Peptide
65	S3AE	179	TYR	Peptide
65	S3AE	181	LEU	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	A16S	32063	0	0	0	0
2	A23S	64339	0	0	0	0
3	AEFG	5677	0	0	0	0
4	A5S	2609	0	1324	47	0
5	AL2P	1754	0	0	0	0
6	AL3P	2695	0	0	0	0
7	AL4P	1926	0	0	0	0
8	AL5P	1343	0	0	0	0
9	AL6P	1431	0	0	0	0
10	ALX0	629	0	0	0	0
11	L10E	1310	0	0	0	0
12	L13P	1109	0	0	0	0
13	L141	669	0	0	0	0
13	L142	669	0	0	0	0
14	L14P	1034	0	0	0	0
15	L15E	1423	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
16	L18E	895	0	0	0	0
17	L18P	1539	0	0	0	0
18	L19E	1206	0	0	0	0
19	L22P	1223	0	0	0	0
20	L23P	650	0	0	0	0
21	L24E	441	0	0	0	0
22	L24P	989	0	0	0	0
23	L29P	513	0	0	0	0
24	L30E	729	0	0	0	0
25	L30P	1254	0	0	0	0
26	L31E	625	0	0	0	0
27	L32E	1010	0	0	0	0
28	L34E	629	0	0	0	0
29	L37A	527	0	0	0	0
30	L37E	436	0	0	0	0
31	L39E	414	0	0	0	0
32	L40E	439	0	0	0	0
33	L44E	753	0	0	0	0
34	L7A1	935	0	0	0	0
34	L7A2	935	0	0	0	0
34	SL7A	935	0	0	0	0
35	L15P	752	0	0	0	0
36	L21E	785	0	0	0	0
37	L45A	816	0	0	0	0
38	L46A	586	0	0	0	0
39	L47A	648	0	0	0	0
40	AL1P	1715	0	0	0	0
41	AS3P	1576	0	0	0	0
42	AS7P	1537	0	0	0	0
43	S10P	824	0	0	0	0
44	S13P	1204	0	0	0	0
45	S14P	432	0	0	0	0
46	S17E	517	0	0	0	0
47	S19E	1230	0	0	0	0
48	S19P	968	0	0	0	0
49	AS9P	1096	0	0	0	0
50	S28E	498	0	0	0	0
51	S27A	435	0	0	0	0
52	AS2P	1587	0	0	0	0
53	AS4E	1925	0	0	0	0
54	AS4P	1370	0	0	0	0
55	AS5P	1600	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	AS6E	805	0	0	0	0
57	AS8E	993	0	0	0	0
58	AS8P	1028	0	0	0	0
59	S11P	960	0	0	0	0
60	S12P	1103	0	0	0	0
61	S15P	1225	0	0	0	0
62	S17P	885	0	0	0	0
63	S24E	759	0	0	0	0
64	S27E	458	0	0	0	0
65	S3AE	1545	0	0	0	0
66	AETN	1619	0	0	0	0
66	APTN	1619	0	0	0	0
67	AMRN	260	0	0	0	0
68	APTP	67	0	0	0	0
69	AEFG	32	0	0	0	0
70	AS2P	171	0	0	0	0
70	AS5P	80	0	0	0	0
70	AS8P	35	0	0	0	0
All	All	173502	0	1324	47	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 22.

All (47) 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)
4:A5S:76:G:H1	4:A5S:106:U:H3	1.32	0.77
4:A5S:17:G:O6	4:A5S:68:C:N4	2.17	0.75
4:A5S:5:C:H2'	4:A5S:6:A:C8	2.34	0.62
4:A5S:41:U:O2'	4:A5S:46:A:N6	2.32	0.61
4:A5S:113:G:O2'	4:A5S:114:C:O5'	2.21	0.59
4:A5S:40:G:N2	4:A5S:45:G:OP2	2.39	0.56
4:A5S:75:G:H2'	4:A5S:76:G:H8	1.73	0.53
4:A5S:121:G:H2'	4:A5S:122:G:H8	1.74	0.53
4:A5S:57:U:H4'	4:A5S:58:A:O5'	2.08	0.52
4:A5S:40:G:H2'	4:A5S:41:U:C6	2.44	0.52
4:A5S:22:C:H2'	4:A5S:23:G:H8	1.76	0.51
4:A5S:66:C:H2'	4:A5S:67:A:C8	2.45	0.51
4:A5S:77:G:H2'	4:A5S:78:G:H8	1.76	0.51
4:A5S:35:G:O2'	4:A5S:36:A:N7	2.43	0.51
4:A5S:22:C:H2'	4:A5S:23:G:C8	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A5S:75:G:H2'	4:A5S:76:G:C8	2.46	0.50
4:A5S:121:G:H2'	4:A5S:122:G:C8	2.47	0.50
4:A5S:89:C:N4	4:A5S:90:G:O6	2.45	0.49
4:A5S:118:G:H2'	4:A5S:119:A:C8	2.48	0.49
4:A5S:95:G:H2'	4:A5S:96:G:H8	1.78	0.49
4:A5S:73:G:H2'	4:A5S:74:A:H8	1.78	0.48
4:A5S:85:G:H2'	4:A5S:86:A:C8	2.49	0.48
4:A5S:35:G:C6	4:A5S:45:G:H1'	2.48	0.48
4:A5S:4:C:HO2'	4:A5S:5:C:H6	1.62	0.47
4:A5S:49:C:H2'	4:A5S:50:C:C6	2.49	0.47
4:A5S:53:A:HO2'	4:A5S:54:A:H8	1.63	0.46
4:A5S:23:G:H2'	4:A5S:24:G:C8	2.51	0.45
4:A5S:95:G:H2'	4:A5S:96:G:C8	2.51	0.45
4:A5S:98:C:H2'	4:A5S:99:C:C6	2.52	0.45
4:A5S:73:G:H2'	4:A5S:74:A:C8	2.51	0.45
4:A5S:3:C:H2'	4:A5S:4:C:C6	2.52	0.45
4:A5S:1:U:H3'	4:A5S:2:G:C8	2.53	0.44
4:A5S:51:G:C2	4:A5S:52:G:H1'	2.52	0.44
4:A5S:19:G:H2'	4:A5S:20:A:H8	1.84	0.43
4:A5S:23:G:H2'	4:A5S:24:G:H8	1.82	0.43
4:A5S:43:U:H2'	4:A5S:44:C:H6	1.83	0.43
4:A5S:32:C:O2'	4:A5S:54:A:N1	2.37	0.42
4:A5S:6:A:N6	4:A5S:119:A:H61	2.17	0.42
4:A5S:4:C:O2'	4:A5S:5:C:O5'	2.37	0.41
4:A5S:35:G:O6	4:A5S:45:G:H1'	2.21	0.41
4:A5S:41:U:HO2'	4:A5S:44:C:N4	2.19	0.41
4:A5S:76:G:C2	4:A5S:107:C:C2	3.09	0.41
4:A5S:52:G:C6	4:A5S:53:A:C6	3.09	0.41
4:A5S:55:G:H2'	4:A5S:56:U:C6	2.56	0.41
4:A5S:40:G:C2	4:A5S:45:G:C2	3.09	0.40
4:A5S:42:U:OP1	4:A5S:43:U:H5''	2.21	0.40
4:A5S:52:G:C2	4:A5S:53:A:C4	3.09	0.40

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	
3	AEFG	721/729 (99%)	614 (85%)	96 (13%)	11 (2%)	8	38
5	AL2P	232/234 (99%)	215 (93%)	17 (7%)	0	100	100
6	AL3P	337/339 (99%)	301 (89%)	36 (11%)	0	100	100
7	AL4P	249/251 (99%)	234 (94%)	15 (6%)	0	100	100
8	AL5P	166/168 (99%)	146 (88%)	20 (12%)	0	100	100
9	AL6P	179/181 (99%)	170 (95%)	9 (5%)	0	100	100
10	ALX0	74/76 (97%)	64 (86%)	10 (14%)	0	100	100
11	L10E	162/164 (99%)	143 (88%)	19 (12%)	0	100	100
12	L13P	138/140 (99%)	122 (88%)	16 (12%)	0	100	100
13	L141	84/86 (98%)	74 (88%)	10 (12%)	0	100	100
13	L142	84/86 (98%)	78 (93%)	6 (7%)	0	100	100
14	L14P	132/134 (98%)	116 (88%)	16 (12%)	0	100	100
15	L15E	167/169 (99%)	159 (95%)	8 (5%)	0	100	100
16	L18E	110/112 (98%)	106 (96%)	4 (4%)	0	100	100
17	L18P	191/193 (99%)	176 (92%)	15 (8%)	0	100	100
18	L19E	142/144 (99%)	139 (98%)	3 (2%)	0	100	100
19	L22P	148/150 (99%)	138 (93%)	10 (7%)	0	100	100
20	L23P	79/81 (98%)	71 (90%)	8 (10%)	0	100	100
21	L24E	52/54 (96%)	48 (92%)	4 (8%)	0	100	100
22	L24P	120/122 (98%)	111 (92%)	9 (8%)	0	100	100
23	L29P	61/63 (97%)	55 (90%)	6 (10%)	0	100	100
24	L30E	92/94 (98%)	89 (97%)	3 (3%)	0	100	100
25	L30P	153/155 (99%)	135 (88%)	18 (12%)	0	100	100
26	L31E	73/75 (97%)	68 (93%)	5 (7%)	0	100	100
27	L32E	121/123 (98%)	110 (91%)	11 (9%)	0	100	100
28	L34E	75/77 (97%)	54 (72%)	20 (27%)	1 (1%)	10	41
29	L37A	63/65 (97%)	56 (89%)	7 (11%)	0	100	100
30	L37E	52/54 (96%)	45 (86%)	7 (14%)	0	100	100
31	L39E	47/49 (96%)	41 (87%)	6 (13%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
32	L40E	53/55 (96%)	35 (66%)	18 (34%)	0	100	100
33	L44E	90/92 (98%)	87 (97%)	3 (3%)	0	100	100
34	L7A1	121/123 (98%)	118 (98%)	3 (2%)	0	100	100
34	L7A2	121/123 (98%)	119 (98%)	2 (2%)	0	100	100
34	SL7A	121/123 (98%)	113 (93%)	8 (7%)	0	100	100
35	L15P	90/144 (62%)	84 (93%)	6 (7%)	0	100	100
36	L21E	95/97 (98%)	89 (94%)	6 (6%)	0	100	100
37	L45A	99/101 (98%)	82 (83%)	14 (14%)	3 (3%)	3	26
38	L46A	68/70 (97%)	64 (94%)	4 (6%)	0	100	100
39	L47A	78/80 (98%)	72 (92%)	6 (8%)	0	100	100
40	AL1P	214/216 (99%)	194 (91%)	20 (9%)	0	100	100
41	AS3P	199/201 (99%)	182 (92%)	17 (8%)	0	100	100
42	AS7P	191/193 (99%)	157 (82%)	34 (18%)	0	100	100
43	S10P	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
44	S13P	145/147 (99%)	130 (90%)	15 (10%)	0	100	100
45	S14P	50/52 (96%)	43 (86%)	7 (14%)	0	100	100
46	S17E	60/62 (97%)	56 (93%)	4 (7%)	0	100	100
47	S19E	145/150 (97%)	134 (92%)	11 (8%)	0	100	100
48	S19P	113/115 (98%)	99 (88%)	14 (12%)	0	100	100
49	AS9P	134/136 (98%)	124 (92%)	10 (8%)	0	100	100
50	S28E	61/63 (97%)	55 (90%)	6 (10%)	0	100	100
51	S27A	52/54 (96%)	40 (77%)	12 (23%)	0	100	100
52	AS2P	194/196 (99%)	181 (93%)	13 (7%)	0	100	100
53	AS4E	238/240 (99%)	208 (87%)	30 (13%)	0	100	100
54	AS4P	164/166 (99%)	134 (82%)	30 (18%)	0	100	100
55	AS5P	202/204 (99%)	182 (90%)	20 (10%)	0	100	100
56	AS6E	103/105 (98%)	84 (82%)	19 (18%)	0	100	100
57	AS8E	124/126 (98%)	109 (88%)	15 (12%)	0	100	100
58	AS8P	128/130 (98%)	122 (95%)	6 (5%)	0	100	100
59	S11P	126/128 (98%)	108 (86%)	18 (14%)	0	100	100
60	S12P	141/143 (99%)	119 (84%)	22 (16%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
61	S15P	147/149 (99%)	131 (89%)	16 (11%)	0	100	100
62	S17P	109/111 (98%)	99 (91%)	10 (9%)	0	100	100
63	S24E	94/96 (98%)	89 (95%)	5 (5%)	0	100	100
64	S27E	57/59 (97%)	52 (91%)	5 (9%)	0	100	100
65	S3AE	187/189 (99%)	156 (83%)	30 (16%)	1 (0%)	25	60
68	APTP	4/6 (67%)	3 (75%)	1 (25%)	0	100	100
All	All	8720/8913 (98%)	7827 (90%)	877 (10%)	16 (0%)	45	75

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AEFG	298	ILE
3	AEFG	356	VAL
3	AEFG	469	ILE
37	L45A	4	VAL
37	L45A	55	VAL
3	AEFG	598	VAL
3	AEFG	659	VAL
3	AEFG	47	ILE
3	AEFG	395	GLU
37	L45A	85	LYS
3	AEFG	71	VAL
3	AEFG	117	VAL
28	L34E	19	PRO
65	S3AE	180	PRO
3	AEFG	25	ILE
3	AEFG	548	ILE

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	
3	AEFG	625/627 (100%)	620 (99%)	5 (1%)	79	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	AL2P	181/181 (100%)	180 (99%)	1 (1%)	84	87
6	AL3P	297/297 (100%)	296 (100%)	1 (0%)	91	91
7	AL4P	212/212 (100%)	210 (99%)	2 (1%)	75	83
8	AL5P	144/144 (100%)	143 (99%)	1 (1%)	81	86
9	AL6P	157/157 (100%)	157 (100%)	0	100	100
10	ALX0	68/68 (100%)	68 (100%)	0	100	100
11	L10E	137/137 (100%)	137 (100%)	0	100	100
12	L13P	121/121 (100%)	119 (98%)	2 (2%)	56	72
13	L141	74/74 (100%)	74 (100%)	0	100	100
13	L142	74/74 (100%)	74 (100%)	0	100	100
14	L14P	110/110 (100%)	109 (99%)	1 (1%)	75	83
15	L15E	146/146 (100%)	145 (99%)	1 (1%)	81	86
16	L18E	98/98 (100%)	96 (98%)	2 (2%)	50	68
17	L18P	162/162 (100%)	160 (99%)	2 (1%)	67	78
18	L19E	126/126 (100%)	126 (100%)	0	100	100
19	L22P	131/131 (100%)	131 (100%)	0	100	100
20	L23P	74/74 (100%)	74 (100%)	0	100	100
21	L24E	50/50 (100%)	50 (100%)	0	100	100
22	L24P	108/108 (100%)	106 (98%)	2 (2%)	52	69
23	L29P	59/59 (100%)	59 (100%)	0	100	100
24	L30E	83/83 (100%)	82 (99%)	1 (1%)	67	78
25	L30P	136/136 (100%)	136 (100%)	0	100	100
26	L31E	66/66 (100%)	66 (100%)	0	100	100
27	L32E	106/106 (100%)	104 (98%)	2 (2%)	52	69
28	L34E	70/70 (100%)	70 (100%)	0	100	100
29	L37A	53/53 (100%)	52 (98%)	1 (2%)	52	69
30	L37E	45/45 (100%)	45 (100%)	0	100	100
31	L39E	44/44 (100%)	43 (98%)	1 (2%)	45	64
32	L40E	50/50 (100%)	48 (96%)	2 (4%)	27	49
33	L44E	84/84 (100%)	84 (100%)	0	100	100
34	L7A1	104/104 (100%)	104 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	L7A2	104/104 (100%)	104 (100%)	0	100	100
34	SL7A	104/104 (100%)	103 (99%)	1 (1%)	73	81
35	L15P	78/118 (66%)	78 (100%)	0	100	100
36	L21E	85/85 (100%)	85 (100%)	0	100	100
37	L45A	91/91 (100%)	89 (98%)	2 (2%)	47	65
38	L46A	66/66 (100%)	65 (98%)	1 (2%)	60	74
39	L47A	74/74 (100%)	73 (99%)	1 (1%)	62	75
40	AL1P	189/190 (100%)	186 (98%)	3 (2%)	58	73
41	AS3P	165/165 (100%)	164 (99%)	1 (1%)	84	87
42	AS7P	166/166 (100%)	163 (98%)	3 (2%)	54	71
43	S10P	92/92 (100%)	89 (97%)	3 (3%)	33	55
44	S13P	129/129 (100%)	129 (100%)	0	100	100
45	S14P	45/45 (100%)	45 (100%)	0	100	100
46	S17E	57/57 (100%)	57 (100%)	0	100	100
47	S19E	133/134 (99%)	132 (99%)	1 (1%)	79	84
48	S19P	106/106 (100%)	105 (99%)	1 (1%)	75	83
49	AS9P	113/113 (100%)	113 (100%)	0	100	100
50	S28E	54/54 (100%)	54 (100%)	0	100	100
51	S27A	47/47 (100%)	47 (100%)	0	100	100
52	AS2P	174/174 (100%)	172 (99%)	2 (1%)	70	79
53	AS4E	210/210 (100%)	207 (99%)	3 (1%)	62	75
54	AS4P	149/149 (100%)	148 (99%)	1 (1%)	81	86
55	AS5P	174/174 (100%)	174 (100%)	0	100	100
56	AS6E	88/88 (100%)	86 (98%)	2 (2%)	45	64
57	AS8E	106/106 (100%)	105 (99%)	1 (1%)	75	83
58	AS8P	111/111 (100%)	110 (99%)	1 (1%)	75	83
59	S11P	94/94 (100%)	94 (100%)	0	100	100
60	S12P	116/116 (100%)	115 (99%)	1 (1%)	75	83
61	S15P	133/133 (100%)	133 (100%)	0	100	100
62	S17P	97/97 (100%)	95 (98%)	2 (2%)	48	66
63	S24E	84/84 (100%)	83 (99%)	1 (1%)	67	78

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
64	S27E	51/51 (100%)	51 (100%)	0	100	100
65	S3AE	170/170 (100%)	168 (99%)	2 (1%)	67	78
68	APTP	6/6 (100%)	6 (100%)	0	100	100
All	All	7656/7700 (99%)	7596 (99%)	60 (1%)	77	84

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	AEFG	262	ARG
3	AEFG	299	ASN
3	AEFG	405	ARG
3	AEFG	408	LYS
3	AEFG	694	ARG
5	AL2P	234	ARG
6	AL3P	4	ARG
7	AL4P	79	ARG
7	AL4P	117	LYS
8	AL5P	54	ARG
12	L13P	33	LYS
12	L13P	147	LYS
14	L14P	37	ILE
15	L15E	25	LYS
16	L18E	36	ARG
16	L18E	117	LYS
17	L18P	138	LYS
17	L18P	140	LYS
22	L24P	35	LYS
22	L24P	71	ARG
24	L30E	56	LYS
27	L32E	6	LYS
27	L32E	7	LYS
29	L37A	21	LYS
31	L39E	18	LYS
32	L40E	15	ARG
32	L40E	32	ARG
37	L45A	87	LYS
37	L45A	98	LYS
38	L46A	67	LYS
39	L47A	31	LYS
40	AL1P	54	ILE
40	AL1P	196	LYS

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Mol	Chain	Res	Type
40	AL1P	204	ARG
41	AS3P	99	ARG
42	AS7P	46	ARG
42	AS7P	66	ASN
42	AS7P	189	ARG
43	S10P	25	ARG
43	S10P	45	ARG
43	S10P	51	MET
47	S19E	118	ARG
48	S19P	58	ARG
34	SL7A	34	LYS
52	AS2P	185	ARG
52	AS2P	204	LYS
53	AS4E	20	LYS
53	AS4E	53	ARG
53	AS4E	196	LYS
54	AS4P	71	LYS
56	AS6E	125	LYS
56	AS6E	213	ARG
57	AS8E	11	LYS
58	AS8P	9	ASN
60	S12P	22	LYS
62	S17P	21	LYS
62	S17P	37	ARG
63	S24E	50	LYS
65	S3AE	17	LYS
65	S3AE	100	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A16S	1491/1494 (99%)	371 (24%)	52 (3%)
2	A23S	2986/3022 (98%)	649 (21%)	37 (1%)
4	A5S	121/122 (99%)	18 (14%)	1 (0%)
66	AETN	75/76 (98%)	31 (41%)	3 (4%)
66	APTN	75/76 (98%)	20 (26%)	1 (1%)
67	AMRN	12/13 (92%)	7 (58%)	0
All	All	4760/4803 (99%)	1096 (23%)	94 (1%)

All (1096) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A16S	9	U
1	A16S	11	C
1	A16S	39	A
1	A16S	41	C
1	A16S	47	A
1	A16S	48	G
1	A16S	49	G
1	A16S	50	G
1	A16S	53	A
1	A16S	54	A
1	A16S	55	G
1	A16S	64	G
1	A16S	67	U
1	A16S	68	U
1	A16S	73	U
1	A16S	77	G
1	A16S	78	G
1	A16S	81	A
1	A16S	82	A
1	A16S	83	G
1	A16S	86	A
1	A16S	103	A
1	A16S	107	A
1	A16S	108	C
1	A16S	109	A
1	A16S	114	G
1	A16S	117	A
1	A16S	118	A
1	A16S	119	C
1	A16S	148	A
1	A16S	149	A
1	A16S	150	A
1	A16S	151	C
1	A16S	159	A
1	A16S	161	U
1	A16S	169	A
1	A16S	185	A
1	A16S	187	U
1	A16S	195	U
1	A16S	200	A
1	A16S	202	A
1	A16S	203	G

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Mol	Chain	Res	Type
1	A16S	205	G
1	A16S	212	G
1	A16S	213	C
1	A16S	214	U
1	A16S	215	A
1	A16S	220	C
1	A16S	221	C
1	A16S	222	G
1	A16S	223	U
1	A16S	230	C
1	A16S	252	C
1	A16S	256	U
1	A16S	257	C
1	A16S	259	G
1	A16S	262	U
1	A16S	263	G
1	A16S	266	G
1	A16S	270	G
1	A16S	274	A
1	A16S	278	G
1	A16S	279	C
1	A16S	294	A
1	A16S	301	G
1	A16S	313	G
1	A16S	320	C
1	A16S	328	U
1	A16S	340	C
1	A16S	341	A
1	A16S	344	G
1	A16S	355	U
1	A16S	356	A
1	A16S	357	C
1	A16S	358	G
1	A16S	359	G
1	A16S	360	G
1	A16S	364	C
1	A16S	365	A
1	A16S	366	C
1	A16S	368	A
1	A16S	377	A
1	A16S	378	C
1	A16S	380	U

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Mol	Chain	Res	Type
1	A16S	381	C
1	A16S	385	A
1	A16S	402	G
1	A16S	404	G
1	A16S	405	C
1	A16S	409	A
1	A16S	418	G
1	A16S	436	U
1	A16S	437	U
1	A16S	438	C
1	A16S	443	C
1	A16S	444	U
1	A16S	445	C
1	A16S	447	A
1	A16S	449	A
1	A16S	451	A
1	A16S	452	G
1	A16S	453	G
1	A16S	459	G
1	A16S	461	A
1	A16S	462	U
1	A16S	463	A
1	A16S	472	G
1	A16S	473	C
1	A16S	474	A
1	A16S	475	A
1	A16S	476	G
1	A16S	477	U
1	A16S	478	C
1	A16S	480	G
1	A16S	482	U
1	A16S	483	G
1	A16S	484	U
1	A16S	485	C
1	A16S	493	G
1	A16S	495	G
1	A16S	496	G
1	A16S	497	U
1	A16S	498	A
1	A16S	499	A
1	A16S	513	A
1	A16S	525	G

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Mol	Chain	Res	Type
1	A16S	526	A
1	A16S	527	U
1	A16S	530	C
1	A16S	538	A
1	A16S	539	A
1	A16S	542	C
1	A16S	543	G
1	A16S	554	G
1	A16S	562	A
1	A16S	585	U
1	A16S	586	C
1	A16S	587	A
1	A16S	588	A
1	A16S	596	A
1	A16S	597	C
1	A16S	598	U
1	A16S	600	G
1	A16S	619	A
1	A16S	620	G
1	A16S	627	G
1	A16S	631	A
1	A16S	652	U
1	A16S	654	G
1	A16S	661	A
1	A16S	667	U
1	A16S	668	A
1	A16S	669	G
1	A16S	685	C
1	A16S	689	U
1	A16S	690	G
1	A16S	697	G
1	A16S	714	G
1	A16S	715	C
1	A16S	721	G
1	A16S	726	G
1	A16S	731	G
1	A16S	732	A
1	A16S	740	G
1	A16S	741	G
1	A16S	743	A
1	A16S	747	A
1	A16S	758	A

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Mol	Chain	Res	Type
1	A16S	759	U
1	A16S	760	A
1	A16S	779	U
1	A16S	783	C
1	A16S	784	G
1	A16S	787	G
1	A16S	793	U
1	A16S	794	A
1	A16S	795	G
1	A16S	810	U
1	A16S	811	A
1	A16S	812	G
1	A16S	813	A
1	A16S	842	A
1	A16S	843	A
1	A16S	859	A
1	A16S	870	A
1	A16S	872	G
1	A16S	884	A
1	A16S	896	G
1	A16S	897	G
1	A16S	904	C
1	A16S	905	C
1	A16S	906	A
1	A16S	909	A
1	A16S	911	G
1	A16S	913	G
1	A16S	930	A
1	A16S	931	U
1	A16S	932	U
1	A16S	933	G
1	A16S	936	G
1	A16S	937	U
1	A16S	938	C
1	A16S	940	A
1	A16S	942	G
1	A16S	945	U
1	A16S	946	G
1	A16S	947	G
1	A16S	949	A
1	A16S	953	U
1	A16S	954	A

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Mol	Chain	Res	Type
1	A16S	959	G
1	A16S	961	G
1	A16S	962	A
1	A16S	963	G
1	A16S	964	A
1	A16S	966	C
1	A16S	971	U
1	A16S	973	U
1	A16S	974	G
1	A16S	975	A
1	A16S	976	C
1	A16S	982	G
1	A16S	985	U
1	A16S	987	A
1	A16S	988	C
1	A16S	989	G
1	A16S	990	A
1	A16S	995	G
1	A16S	998	U
1	A16S	999	G
1	A16S	1000	A
1	A16S	1001	C
1	A16S	1005	C
1	A16S	1010	A
1	A16S	1013	A
1	A16S	1014	G
1	A16S	1017	G
1	A16S	1025	G
1	A16S	1028	G
1	A16S	1030	C
1	A16S	1031	A
1	A16S	1032	G
1	A16S	1044	A
1	A16S	1045	A
1	A16S	1047	U
1	A16S	1048	G
1	A16S	1049	U
1	A16S	1050	C
1	A16S	1055	U
1	A16S	1058	G
1	A16S	1059	U
1	A16S	1061	C

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Mol	Chain	Res	Type
1	A16S	1062	G
1	A16S	1065	A
1	A16S	1066	A
1	A16S	1088	G
1	A16S	1089	U
1	A16S	1090	U
1	A16S	1091	G
1	A16S	1093	U
1	A16S	1094	A
1	A16S	1095	U
1	A16S	1098	U
1	A16S	1099	G
1	A16S	1102	C
1	A16S	1104	C
1	A16S	1107	G
1	A16S	1109	C
1	A16S	1112	G
1	A16S	1115	C
1	A16S	1117	A
1	A16S	1120	C
1	A16S	1123	G
1	A16S	1128	A
1	A16S	1129	C
1	A16S	1130	U
1	A16S	1134	G
1	A16S	1135	G
1	A16S	1137	G
1	A16S	1138	U
1	A16S	1139	A
1	A16S	1152	G
1	A16S	1153	A
1	A16S	1154	G
1	A16S	1166	A
1	A16S	1167	G
1	A16S	1170	C
1	A16S	1172	G
1	A16S	1181	G
1	A16S	1182	A
1	A16S	1184	A
1	A16S	1185	C
1	A16S	1186	U
1	A16S	1188	C

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Mol	Chain	Res	Type
1	A16S	1192	G
1	A16S	1195	G
1	A16S	1196	C
1	A16S	1210	U
1	A16S	1211	G
1	A16S	1227	U
1	A16S	1228	G
1	A16S	1230	U
1	A16S	1236	G
1	A16S	1237	A
1	A16S	1238	A
1	A16S	1240	G
1	A16S	1245	A
1	A16S	1246	G
1	A16S	1247	C
1	A16S	1248	C
1	A16S	1249	A
1	A16S	1251	U
1	A16S	1252	C
1	A16S	1253	C
1	A16S	1254	U
1	A16S	1255	U
1	A16S	1259	C
1	A16S	1263	G
1	A16S	1266	G
1	A16S	1267	C
1	A16S	1270	U
1	A16S	1271	U
1	A16S	1274	G
1	A16S	1289	C
1	A16S	1290	C
1	A16S	1291	C
1	A16S	1292	G
1	A16S	1297	C
1	A16S	1306	G
1	A16S	1315	A
1	A16S	1316	G
1	A16S	1317	U
1	A16S	1322	G
1	A16S	1323	C
1	A16S	1327	U
1	A16S	1328	C

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Mol	Chain	Res	Type
1	A16S	1329	A
1	A16S	1330	A
1	A16S	1332	A
1	A16S	1333	A
1	A16S	1334	C
1	A16S	1342	G
1	A16S	1344	A
1	A16S	1347	C
1	A16S	1348	G
1	A16S	1350	C
1	A16S	1355	C
1	A16S	1363	A
1	A16S	1364	C
1	A16S	1366	C
1	A16S	1367	A
1	A16S	1368	C
1	A16S	1369	C
1	A16S	1370	G
1	A16S	1382	A
1	A16S	1388	G
1	A16S	1414	A
1	A16S	1416	A
1	A16S	1430	A
1	A16S	1443	G
1	A16S	1454	A
1	A16S	1460	A
1	A16S	1463	A
1	A16S	1465	G
1	A16S	1467	U
1	A16S	1479	A
1	A16S	1481	C
1	A16S	1490	G
1	A16S	1493	U
1	A16S	1495	A
1	A16S	1497	C
2	A23S	7	G
2	A23S	16	U
2	A23S	17	G
2	A23S	18	G
2	A23S	19	A
2	A23S	25	C
2	A23S	30	C

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Mol	Chain	Res	Type
2	A23S	31	G
2	A23S	36	C
2	A23S	40	A
2	A23S	41	G
2	A23S	53	A
2	A23S	54	A
2	A23S	61	A
2	A23S	64	U
2	A23S	65	G
2	A23S	75	G
2	A23S	80	A
2	A23S	81	A
2	A23S	90	U
2	A23S	91	G
2	A23S	107	A
2	A23S	109	U
2	A23S	110	G
2	A23S	114	U
2	A23S	115	A
2	A23S	116	U
2	A23S	117	C
2	A23S	126	U
2	A23S	137	C
2	A23S	144	U
2	A23S	145	A
2	A23S	147	A
2	A23S	160	A
2	A23S	175	A
2	A23S	179	U
2	A23S	194	G
2	A23S	195	A
2	A23S	200	A
2	A23S	201	A
2	A23S	204	C
2	A23S	206	A
2	A23S	207	U
2	A23S	208	U
2	A23S	209	G
2	A23S	212	A
2	A23S	222	U
2	A23S	227	G
2	A23S	229	G

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Mol	Chain	Res	Type
2	A23S	242	A
2	A23S	243	C
2	A23S	244	A
2	A23S	245	G
2	A23S	246	C
2	A23S	247	C
2	A23S	257	C
2	A23S	268	U
2	A23S	269	A
2	A23S	271	G
2	A23S	281	A
2	A23S	283	G
2	A23S	288	G
2	A23S	289	U
2	A23S	292	C
2	A23S	293	G
2	A23S	304	U
2	A23S	317	C
2	A23S	318	G
2	A23S	319	G
2	A23S	326	U
2	A23S	335	U
2	A23S	336	A
2	A23S	337	G
2	A23S	344	U
2	A23S	345	C
2	A23S	367	A
2	A23S	372	G
2	A23S	373	A
2	A23S	376	G
2	A23S	383	A
2	A23S	385	G
2	A23S	391	G
2	A23S	392	A
2	A23S	403	G
2	A23S	404	U
2	A23S	405	G
2	A23S	418	G
2	A23S	419	U
2	A23S	420	A
2	A23S	433	U
2	A23S	434	U

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Mol	Chain	Res	Type
2	A23S	444	A
2	A23S	448	U
2	A23S	454	A
2	A23S	455	C
2	A23S	460	G
2	A23S	468	G
2	A23S	469	G
2	A23S	470	C
2	A23S	472	A
2	A23S	476	A
2	A23S	494	C
2	A23S	495	A
2	A23S	498	C
2	A23S	499	U
2	A23S	500	A
2	A23S	501	A
2	A23S	502	G
2	A23S	511	G
2	A23S	517	G
2	A23S	525	G
2	A23S	534	A
2	A23S	535	A
2	A23S	536	G
2	A23S	547	A
2	A23S	548	A
2	A23S	549	G
2	A23S	551	G
2	A23S	552	C
2	A23S	553	C
2	A23S	573	A
2	A23S	574	G
2	A23S	575	G
2	A23S	576	G
2	A23S	585	G
2	A23S	586	A
2	A23S	587	A
2	A23S	588	A
2	A23S	589	G
2	A23S	592	G
2	A23S	594	G
2	A23S	621	G
2	A23S	628	G

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Mol	Chain	Res	Type
2	A23S	638	G
2	A23S	639	G
2	A23S	640	G
2	A23S	657	U
2	A23S	658	U
2	A23S	659	U
2	A23S	660	C
2	A23S	667	A
2	A23S	668	A
2	A23S	669	A
2	A23S	670	C
2	A23S	671	A
2	A23S	674	G
2	A23S	675	G
2	A23S	683	G
2	A23S	699	A
2	A23S	720	A
2	A23S	727	A
2	A23S	737	A
2	A23S	740	G
2	A23S	741	C
2	A23S	743	C
2	A23S	753	U
2	A23S	754	U
2	A23S	755	U
2	A23S	757	U
2	A23S	758	A
2	A23S	760	G
2	A23S	762	G
2	A23S	768	C
2	A23S	769	A
2	A23S	778	C
2	A23S	780	G
2	A23S	787	A
2	A23S	801	A
2	A23S	802	G
2	A23S	808	A
2	A23S	810	A
2	A23S	819	U
2	A23S	828	G
2	A23S	829	G
2	A23S	838	A

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Mol	Chain	Res	Type
2	A23S	844	G
2	A23S	847	G
2	A23S	848	A
2	A23S	849	A
2	A23S	850	A
2	A23S	851	C
2	A23S	857	U
2	A23S	858	G
2	A23S	859	G
2	A23S	863	G
2	A23S	868	A
2	A23S	869	U
2	A23S	875	U
2	A23S	878	U
2	A23S	879	G
2	A23S	883	U
2	A23S	884	G
2	A23S	892	U
2	A23S	911	G
2	A23S	912	G
2	A23S	915	C
2	A23S	918	A
2	A23S	920	G
2	A23S	921	A
2	A23S	925	A
2	A23S	928	U
2	A23S	936	U
2	A23S	937	G
2	A23S	941	G
2	A23S	947	U
2	A23S	948	C
2	A23S	957	A
2	A23S	963	C
2	A23S	964	C
2	A23S	966	A
2	A23S	981	A
2	A23S	982	A
2	A23S	986	A
2	A23S	996	G
2	A23S	1001	G
2	A23S	1004	A
2	A23S	1046	A

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Mol	Chain	Res	Type
2	A23S	1048	C
2	A23S	1053	A
2	A23S	1055	C
2	A23S	1068	A
2	A23S	1069	A
2	A23S	1070	G
2	A23S	1072	A
2	A23S	1073	G
2	A23S	1084	G
2	A23S	1091	G
2	A23S	1096	A
2	A23S	1098	G
2	A23S	1099	G
2	A23S	1102	G
2	A23S	1110	A
2	A23S	1111	A
2	A23S	1122	C
2	A23S	1126	G
2	A23S	1127	A
2	A23S	1145	A
2	A23S	1148	U
2	A23S	1149	C
2	A23S	1158	G
2	A23S	1163	A
2	A23S	1165	C
2	A23S	1166	G
2	A23S	1168	A
2	A23S	1173	C
2	A23S	1174	G
2	A23S	1177	U
2	A23S	1178	C
2	A23S	1179	C
2	A23S	1180	A
2	A23S	1185	U
2	A23S	1187	G
2	A23S	1188	A
2	A23S	1196	A
2	A23S	1198	G
2	A23S	1199	G
2	A23S	1200	U
2	A23S	1201	G
2	A23S	1202	G

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Mol	Chain	Res	Type
2	A23S	1203	G
2	A23S	1205	C
2	A23S	1206	C
2	A23S	1207	A
2	A23S	1208	G
2	A23S	1211	G
2	A23S	1213	A
2	A23S	1217	A
2	A23S	1219	C
2	A23S	1221	U
2	A23S	1222	C
2	A23S	1225	A
2	A23S	1227	G
2	A23S	1228	A
2	A23S	1229	G
2	A23S	1230	U
2	A23S	1233	G
2	A23S	1237	C
2	A23S	1244	C
2	A23S	1245	C
2	A23S	1249	C
2	A23S	1251	A
2	A23S	1252	G
2	A23S	1269	A
2	A23S	1270	U
2	A23S	1271	U
2	A23S	1272	G
2	A23S	1273	G
2	A23S	1275	C
2	A23S	1282	C
2	A23S	1283	A
2	A23S	1289	G
2	A23S	1296	A
2	A23S	1317	C
2	A23S	1318	U
2	A23S	1319	A
2	A23S	1320	G
2	A23S	1322	G
2	A23S	1328	C
2	A23S	1330	G
2	A23S	1339	G
2	A23S	1344	G

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Mol	Chain	Res	Type
2	A23S	1348	U
2	A23S	1350	G
2	A23S	1351	G
2	A23S	1354	A
2	A23S	1355	G
2	A23S	1368	G
2	A23S	1370	G
2	A23S	1379	G
2	A23S	1381	C
2	A23S	1383	C
2	A23S	1391	G
2	A23S	1393	G
2	A23S	1395	A
2	A23S	1396	G
2	A23S	1398	U
2	A23S	1399	C
2	A23S	1406	G
2	A23S	1408	A
2	A23S	1417	G
2	A23S	1424	G
2	A23S	1428	A
2	A23S	1440	G
2	A23S	1444	G
2	A23S	1445	A
2	A23S	1458	U
2	A23S	1465	A
2	A23S	1482	A
2	A23S	1485	U
2	A23S	1489	C
2	A23S	1490	G
2	A23S	1501	A
2	A23S	1504	G
2	A23S	1506	C
2	A23S	1515	A
2	A23S	1525	A
2	A23S	1529	A
2	A23S	1530	A
2	A23S	1540	C
2	A23S	1541	A
2	A23S	1542	U
2	A23S	1543	U
2	A23S	1545	C

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Mol	Chain	Res	Type
2	A23S	1549	G
2	A23S	1555	C
2	A23S	1564	G
2	A23S	1565	C
2	A23S	1569	A
2	A23S	1571	C
2	A23S	1574	A
2	A23S	1575	A
2	A23S	1576	G
2	A23S	1579	A
2	A23S	1581	A
2	A23S	1587	G
2	A23S	1589	C
2	A23S	1598	G
2	A23S	1601	G
2	A23S	1613	C
2	A23S	1615	A
2	A23S	1617	C
2	A23S	1618	G
2	A23S	1619	U
2	A23S	1625	C
2	A23S	1626	C
2	A23S	1633	U
2	A23S	1634	C
2	A23S	1635	A
2	A23S	1636	A
2	A23S	1644	G
2	A23S	1645	A
2	A23S	1646	G
2	A23S	1647	A
2	A23S	1648	G
2	A23S	1659	G
2	A23S	1661	G
2	A23S	1666	G
2	A23S	1667	G
2	A23S	1671	A
2	A23S	1676	G
2	A23S	1683	G
2	A23S	1684	C
2	A23S	1685	C
2	A23S	1690	G
2	A23S	1692	U

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Mol	Chain	Res	Type
2	A23S	1693	A
2	A23S	1694	G
2	A23S	1701	U
2	A23S	1707	G
2	A23S	1708	A
2	A23S	1713	U
2	A23S	1715	G
2	A23S	1720	C
2	A23S	1721	A
2	A23S	1724	A
2	A23S	1729	G
2	A23S	1740	C
2	A23S	1741	G
2	A23S	1742	A
2	A23S	1743	U
2	A23S	1745	C
2	A23S	1751	G
2	A23S	1752	A
2	A23S	1756	U
2	A23S	1762	G
2	A23S	1764	A
2	A23S	1770	A
2	A23S	1771	C
2	A23S	1772	U
2	A23S	1776	G
2	A23S	1786	G
2	A23S	1789	A
2	A23S	1801	U
2	A23S	1804	G
2	A23S	1809	G
2	A23S	1810	U
2	A23S	1811	A
2	A23S	1850	U
2	A23S	1857	A
2	A23S	1868	A
2	A23S	1869	U
2	A23S	1877	U
2	A23S	1879	A
2	A23S	1882	A
2	A23S	1885	C
2	A23S	1900	G
2	A23S	1907	G

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Mol	Chain	Res	Type
2	A23S	1908	A
2	A23S	1917	A
2	A23S	1928	A
2	A23S	1935	A
2	A23S	1953	A
2	A23S	1955	G
2	A23S	1965	G
2	A23S	1972	A
2	A23S	1973	A
2	A23S	1979	G
2	A23S	2007	U
2	A23S	2008	C
2	A23S	2010	A
2	A23S	2013	G
2	A23S	2020	G
2	A23S	2028	G
2	A23S	2030	A
2	A23S	2032	G
2	A23S	2036	G
2	A23S	2041	A
2	A23S	2043	C
2	A23S	2044	U
2	A23S	2045	C
2	A23S	2046	U
2	A23S	2047	G
2	A23S	2054	U
2	A23S	2055	U
2	A23S	2056	A
2	A23S	2058	G
2	A23S	2059	G
2	A23S	2060	U
2	A23S	2066	A
2	A23S	2081	A
2	A23S	2084	U
2	A23S	2092	C
2	A23S	2093	G
2	A23S	2094	C
2	A23S	2096	U
2	A23S	2099	A
2	A23S	2100	U
2	A23S	2101	G
2	A23S	2105	C

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Mol	Chain	Res	Type
2	A23S	2120	U
2	A23S	2122	U
2	A23S	2123	C
2	A23S	2125	C
2	A23S	2149	U
2	A23S	2150	G
2	A23S	2151	A
2	A23S	2159	C
2	A23S	2160	A
2	A23S	2161	C
2	A23S	2162	A
2	A23S	2163	G
2	A23S	2170	A
2	A23S	2180	A
2	A23S	2181	C
2	A23S	2184	A
2	A23S	2189	A
2	A23S	2190	G
2	A23S	2191	A
2	A23S	2198	G
2	A23S	2209	G
2	A23S	2222	G
2	A23S	2225	C
2	A23S	2231	G
2	A23S	2234	U
2	A23S	2237	A
2	A23S	2239	G
2	A23S	2240	C
2	A23S	2241	G
2	A23S	2244	G
2	A23S	2245	A
2	A23S	2247	U
2	A23S	2248	A
2	A23S	2249	G
2	A23S	2252	G
2	A23S	2255	A
2	A23S	2260	U
2	A23S	2261	C
2	A23S	2262	G
2	A23S	2265	C
2	A23S	2266	C
2	A23S	2268	G

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Mol	Chain	Res	Type
2	A23S	2270	C
2	A23S	2274	U
2	A23S	2276	G
2	A23S	2279	G
2	A23S	2285	G
2	A23S	2287	A
2	A23S	2288	C
2	A23S	2290	C
2	A23S	2292	A
2	A23S	2293	A
2	A23S	2298	A
2	A23S	2299	A
2	A23S	2300	A
2	A23S	2302	A
2	A23S	2319	G
2	A23S	2320	A
2	A23S	2327	A
2	A23S	2328	A
2	A23S	2337	A
2	A23S	2346	A
2	A23S	2359	G
2	A23S	2360	G
2	A23S	2367	G
2	A23S	2379	A
2	A23S	2382	C
2	A23S	2389	A
2	A23S	2397	C
2	A23S	2401	G
2	A23S	2405	C
2	A23S	2409	A
2	A23S	2410	A
2	A23S	2427	A
2	A23S	2428	C
2	A23S	2429	A
2	A23S	2431	A
2	A23S	2434	G
2	A23S	2441	U
2	A23S	2442	A
2	A23S	2444	A
2	A23S	2449	A
2	A23S	2456	A
2	A23S	2457	A

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Mol	Chain	Res	Type
2	A23S	2465	C
2	A23S	2467	G
2	A23S	2479	C
2	A23S	2500	G
2	A23S	2501	A
2	A23S	2504	G
2	A23S	2505	C
2	A23S	2507	G
2	A23S	2508	G
2	A23S	2510	C
2	A23S	2517	A
2	A23S	2526	G
2	A23S	2527	C
2	A23S	2531	C
2	A23S	2532	A
2	A23S	2540	G
2	A23S	2551	G
2	A23S	2552	A
2	A23S	2555	G
2	A23S	2558	A
2	A23S	2565	C
2	A23S	2571	G
2	A23S	2572	A
2	A23S	2576	C
2	A23S	2596	G
2	A23S	2597	C
2	A23S	2599	C
2	A23S	2600	A
2	A23S	2605	G
2	A23S	2618	G
2	A23S	2620	U
2	A23S	2621	A
2	A23S	2622	C
2	A23S	2626	G
2	A23S	2627	A
2	A23S	2642	A
2	A23S	2653	G
2	A23S	2671	A
2	A23S	2678	C
2	A23S	2687	U
2	A23S	2688	A
2	A23S	2690	A

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Mol	Chain	Res	Type
2	A23S	2691	G
2	A23S	2696	G
2	A23S	2702	G
2	A23S	2705	G
2	A23S	2726	A
2	A23S	2737	U
2	A23S	2745	G
2	A23S	2754	G
2	A23S	2761	U
2	A23S	2762	G
2	A23S	2763	A
2	A23S	2764	G
2	A23S	2769	A
2	A23S	2770	G
2	A23S	2814	U
2	A23S	2826	G
2	A23S	2835	A
2	A23S	2836	U
2	A23S	2837	G
2	A23S	2849	C
2	A23S	2855	G
2	A23S	2870	A
2	A23S	2872	A
2	A23S	2879	A
2	A23S	2886	A
2	A23S	2887	A
2	A23S	2888	C
2	A23S	2891	C
2	A23S	2893	C
2	A23S	2896	C
2	A23S	2898	A
2	A23S	2899	A
2	A23S	2900	A
2	A23S	2924	G
2	A23S	2934	U
2	A23S	2939	G
2	A23S	2940	A
2	A23S	2947	U
2	A23S	2949	A
2	A23S	2952	G
2	A23S	2959	G
2	A23S	2961	U

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Mol	Chain	Res	Type
2	A23S	2962	G
2	A23S	2973	G
2	A23S	2983	C
2	A23S	2994	A
2	A23S	2999	G
2	A23S	3000	C
2	A23S	3005	C
2	A23S	3006	C
2	A23S	3010	U
2	A23S	3019	C
2	A23S	3020	C
2	A23S	3021	C
4	A5S	5	C
4	A5S	10	G
4	A5S	16	A
4	A5S	25	G
4	A5S	27	A
4	A5S	42	U
4	A5S	43	U
4	A5S	45	G
4	A5S	46	A
4	A5S	52	G
4	A5S	53	A
4	A5S	58	A
4	A5S	67	A
4	A5S	68	C
4	A5S	69	G
4	A5S	113	G
4	A5S	114	C
4	A5S	119	A
66	APTN	7	A
66	APTN	8	U
66	APTN	17	C
66	APTN	18	G
66	APTN	20	U
66	APTN	21	A
66	APTN	22	G
66	APTN	40	C
66	APTN	48	C
66	APTN	49	C
66	APTN	50	U
66	APTN	52	G

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Mol	Chain	Res	Type
66	APTN	58	A
66	APTN	61	C
66	APTN	64	A
66	APTN	66	U
66	APTN	67	C
66	APTN	68	C
66	APTN	74	C
66	APTN	76	A
66	AETN	2	C
66	AETN	7	A
66	AETN	8	U
66	AETN	10	G
66	AETN	13	C
66	AETN	16	U
66	AETN	17	C
66	AETN	19	G
66	AETN	20	U
66	AETN	21	A
66	AETN	26	A
66	AETN	31	A
66	AETN	44	G
66	AETN	45	U
66	AETN	46	G
66	AETN	47	U
66	AETN	48	C
66	AETN	49	C
66	AETN	50	U
66	AETN	52	G
66	AETN	53	G
66	AETN	54	U
66	AETN	55	U
66	AETN	56	C
66	AETN	58	A
66	AETN	59	U
66	AETN	65	G
66	AETN	66	U
66	AETN	72	C
66	AETN	73	A
66	AETN	76	A
67	AMRN	-3	U
67	AMRN	-2	U
67	AMRN	-1	U

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Mol	Chain	Res	Type
67	AMRN	1	U
67	AMRN	4	U
67	AMRN	5	U
67	AMRN	7	U

All (94) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A16S	47	A
1	A16S	49	G
1	A16S	54	A
1	A16S	63	A
1	A16S	102	G
1	A16S	117	A
1	A16S	186	A
1	A16S	220	C
1	A16S	255	A
1	A16S	262	U
1	A16S	278	G
1	A16S	293	A
1	A16S	340	C
1	A16S	364	C
1	A16S	384	C
1	A16S	444	U
1	A16S	471	G
1	A16S	475	A
1	A16S	484	U
1	A16S	541	G
1	A16S	653	A
1	A16S	684	A
1	A16S	758	A
1	A16S	778	G
1	A16S	883	A
1	A16S	962	A
1	A16S	972	A
1	A16S	973	U
1	A16S	975	A
1	A16S	987	A
1	A16S	989	G
1	A16S	997	C
1	A16S	999	G
1	A16S	1013	A

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Mol	Chain	Res	Type
1	A16S	1065	A
1	A16S	1116	C
1	A16S	1138	U
1	A16S	1153	A
1	A16S	1171	A
1	A16S	1227	U
1	A16S	1254	U
1	A16S	1269	G
1	A16S	1270	U
1	A16S	1315	A
1	A16S	1316	G
1	A16S	1329	A
1	A16S	1332	A
1	A16S	1349	U
1	A16S	1429	A
1	A16S	1459	U
1	A16S	1489	U
1	A16S	1492	A
2	A23S	136	U
2	A23S	146	U
2	A23S	280	G
2	A23S	336	A
2	A23S	382	U
2	A23S	402	A
2	A23S	404	U
2	A23S	547	A
2	A23S	574	G
2	A23S	668	A
2	A23S	740	G
2	A23S	843	G
2	A23S	920	G
2	A23S	1003	G
2	A23S	1177	U
2	A23S	1179	C
2	A23S	1198	G
2	A23S	1205	C
2	A23S	1224	A
2	A23S	1250	G
2	A23S	1319	A
2	A23S	1327	U
2	A23S	1500	U
2	A23S	1646	G

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Mol	Chain	Res	Type
2	A23S	1647	A
2	A23S	1689	C
2	A23S	1706	U
2	A23S	1741	G
2	A23S	1750	A
2	A23S	2148	C
2	A23S	2504	G
2	A23S	2598	U
2	A23S	2704	U
2	A23S	2725	C
2	A23S	2835	A
2	A23S	2878	U
2	A23S	2923	G
4	A5S	57	U
66	APTn	19	G
66	AETn	47	U
66	AETn	65	G
66	AETn	75	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 58 ligands modelled in this entry, 57 are unknown - leaving 1 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
69	GNP	AEFG	801	-	29,34,34	1.53	7 (24%)	33,54,54	2.13	6 (18%)

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
69	GNP	AEFG	801	-	-	2/14/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
69	AEFG	801	GNP	PB-O3A	3.87	1.63	1.59
69	AEFG	801	GNP	C6-N1	3.44	1.39	1.33
69	AEFG	801	GNP	PG-N3B	2.83	1.70	1.63
69	AEFG	801	GNP	PB-O1B	2.76	1.50	1.46
69	AEFG	801	GNP	PG-O1G	2.51	1.50	1.46
69	AEFG	801	GNP	PB-O2B	-2.40	1.50	1.56
69	AEFG	801	GNP	C5-C6	2.16	1.45	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
69	AEFG	801	GNP	C5-C6-N1	-8.31	112.07	123.43
69	AEFG	801	GNP	C2-N1-C6	5.71	125.01	115.93
69	AEFG	801	GNP	C4-C5-C6	-3.13	117.81	120.80
69	AEFG	801	GNP	PB-O3A-PA	-2.92	122.35	132.62
69	AEFG	801	GNP	N3-C2-N1	-2.80	123.48	127.22
69	AEFG	801	GNP	C2-N3-C4	-2.09	112.97	115.36

There are no chirality outliers.

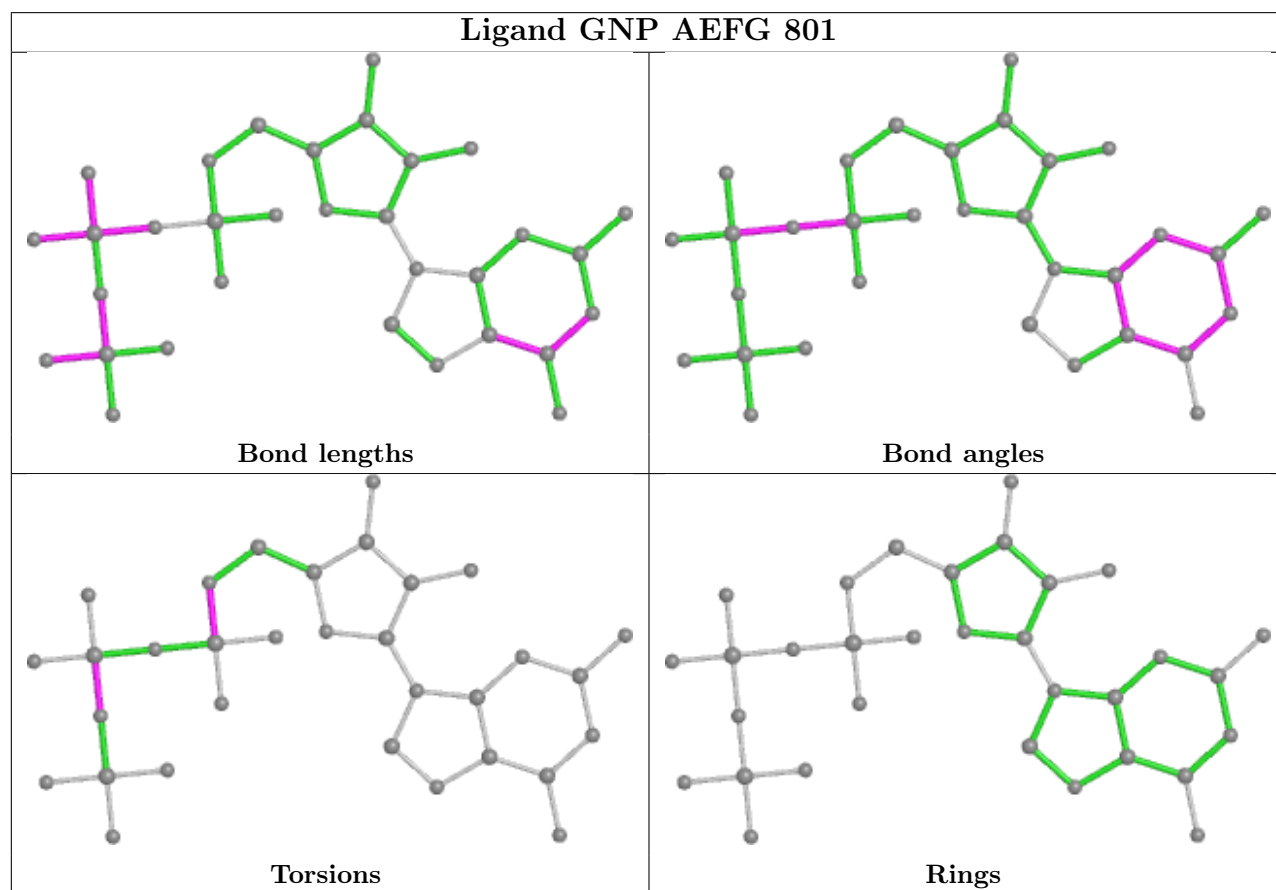
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
69	AEFG	801	GNP	PG-N3B-PB-O1B
69	AEFG	801	GNP	C5'-O5'-PA-O1A

There are no ring outliers.

No monomer is involved in short contacts.

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

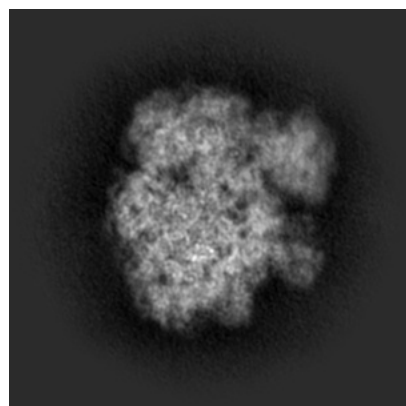
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34866. These allow visual inspection of the internal detail of the map and identification of artifacts.

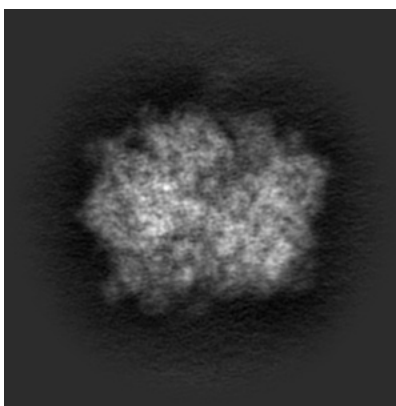
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

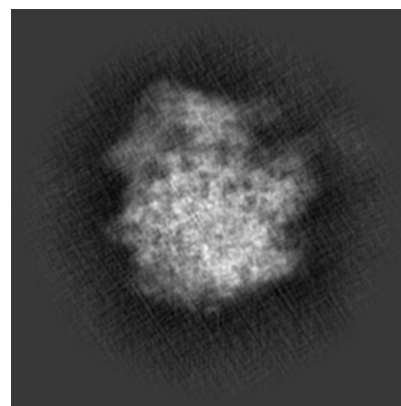
6.1.1 Primary map



X

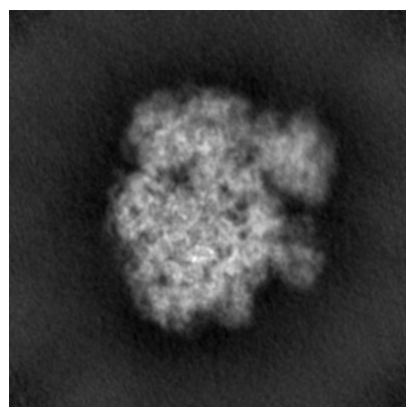


Y

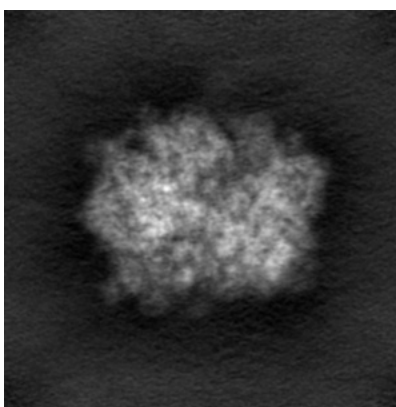


Z

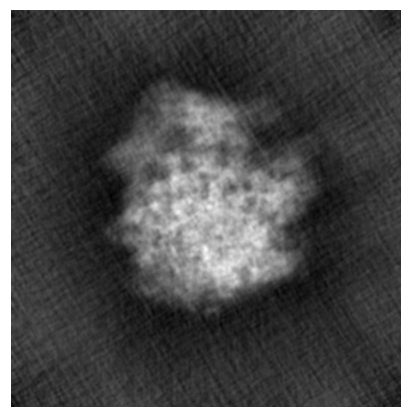
6.1.2 Raw map



X



Y

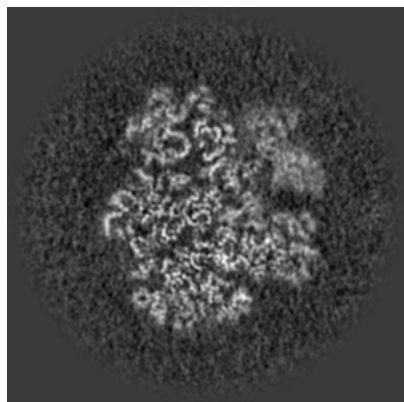


Z

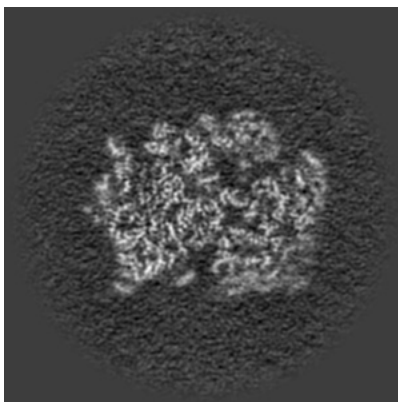
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

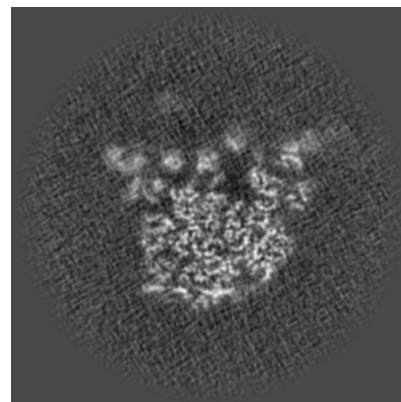
6.2.1 Primary map



X Index: 190

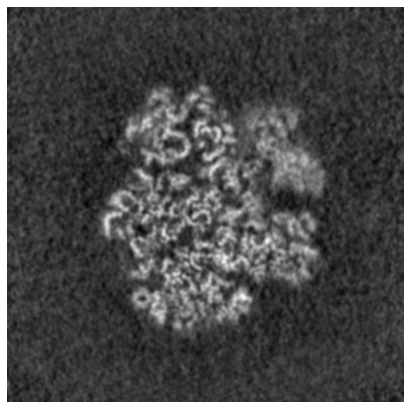


Y Index: 190

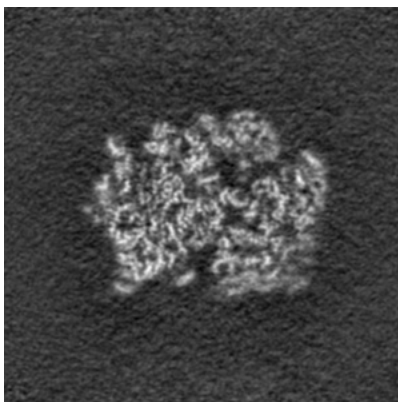


Z Index: 190

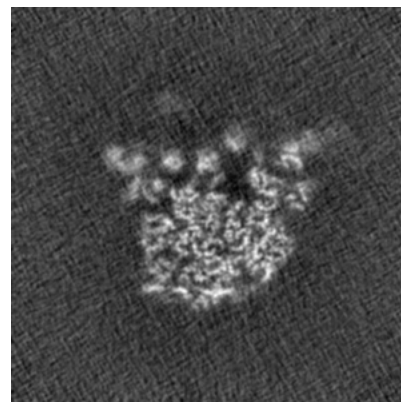
6.2.2 Raw map



X Index: 190



Y Index: 190

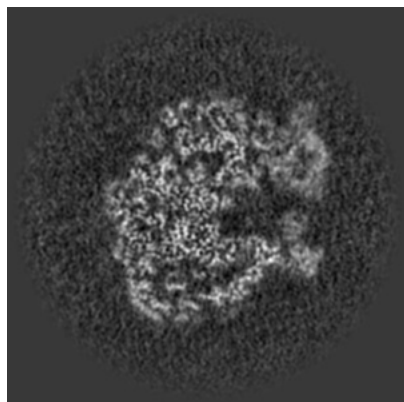


Z Index: 190

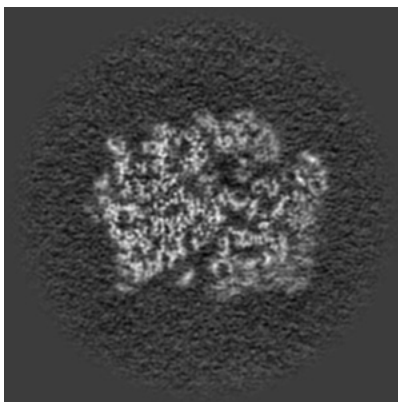
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

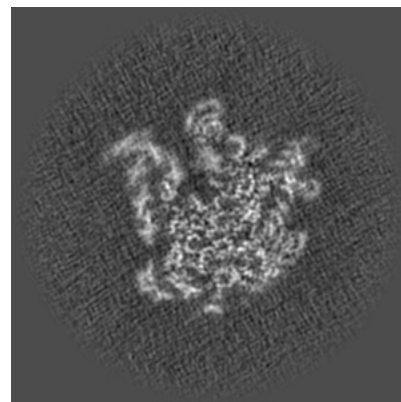
6.3.1 Primary map



X Index: 171

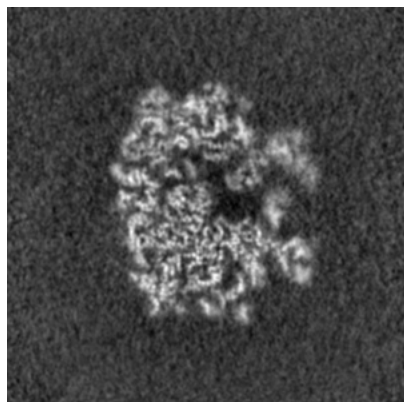


Y Index: 192

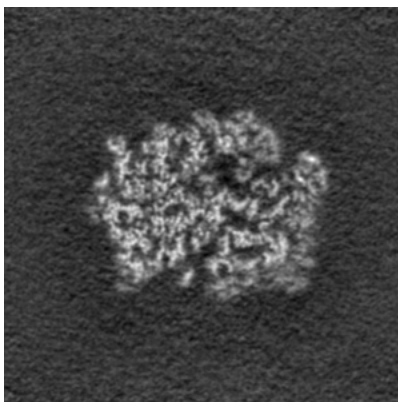


Z Index: 173

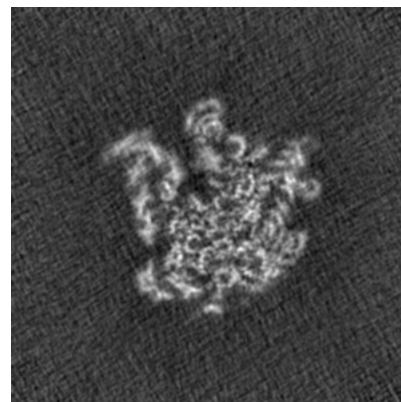
6.3.2 Raw map



X Index: 209



Y Index: 193

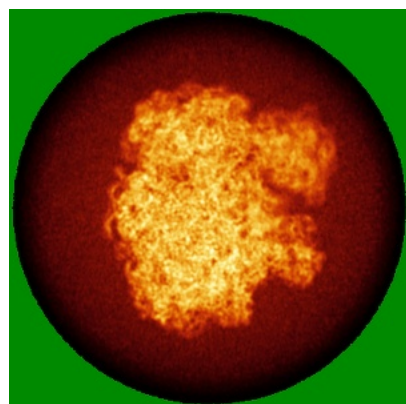


Z Index: 173

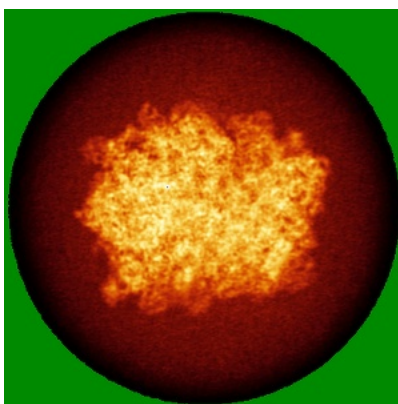
The images above show the largest variance slices of the map in three orthogonal directions.

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

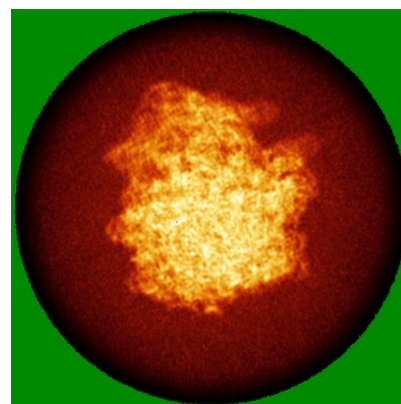
6.4.1 Primary map



X

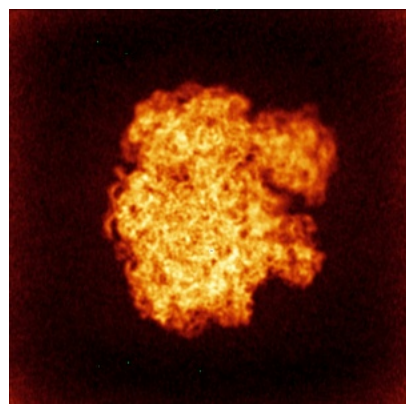


Y

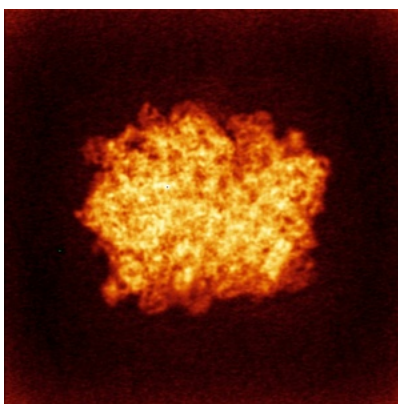


Z

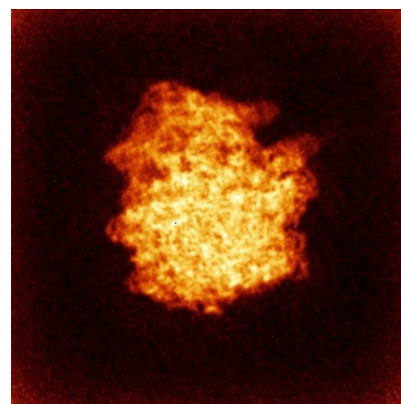
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

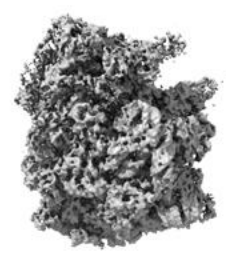
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.33. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

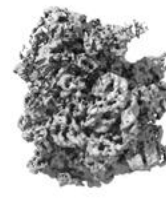
6.5.2 Raw map



X



Y



Z

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

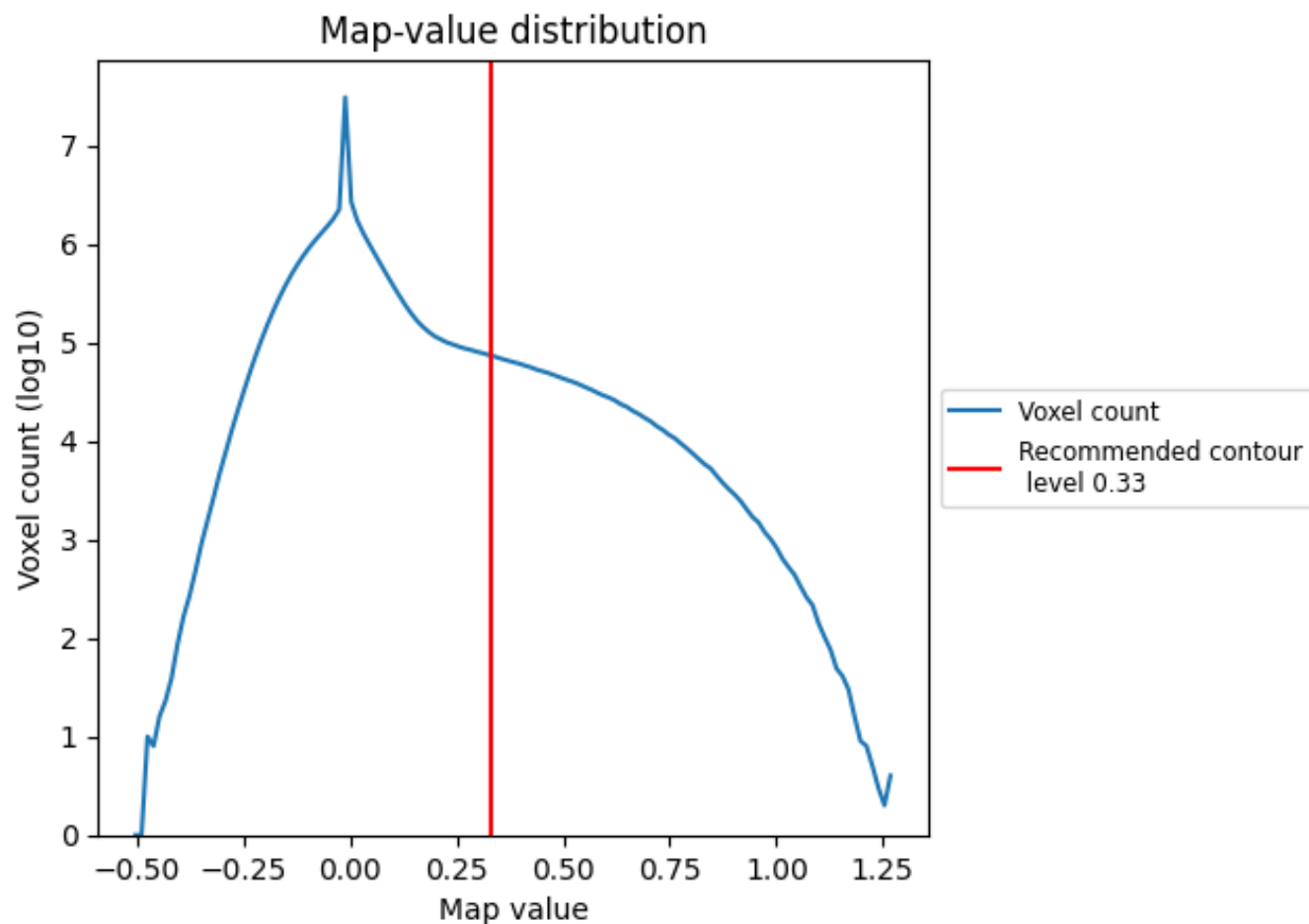
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

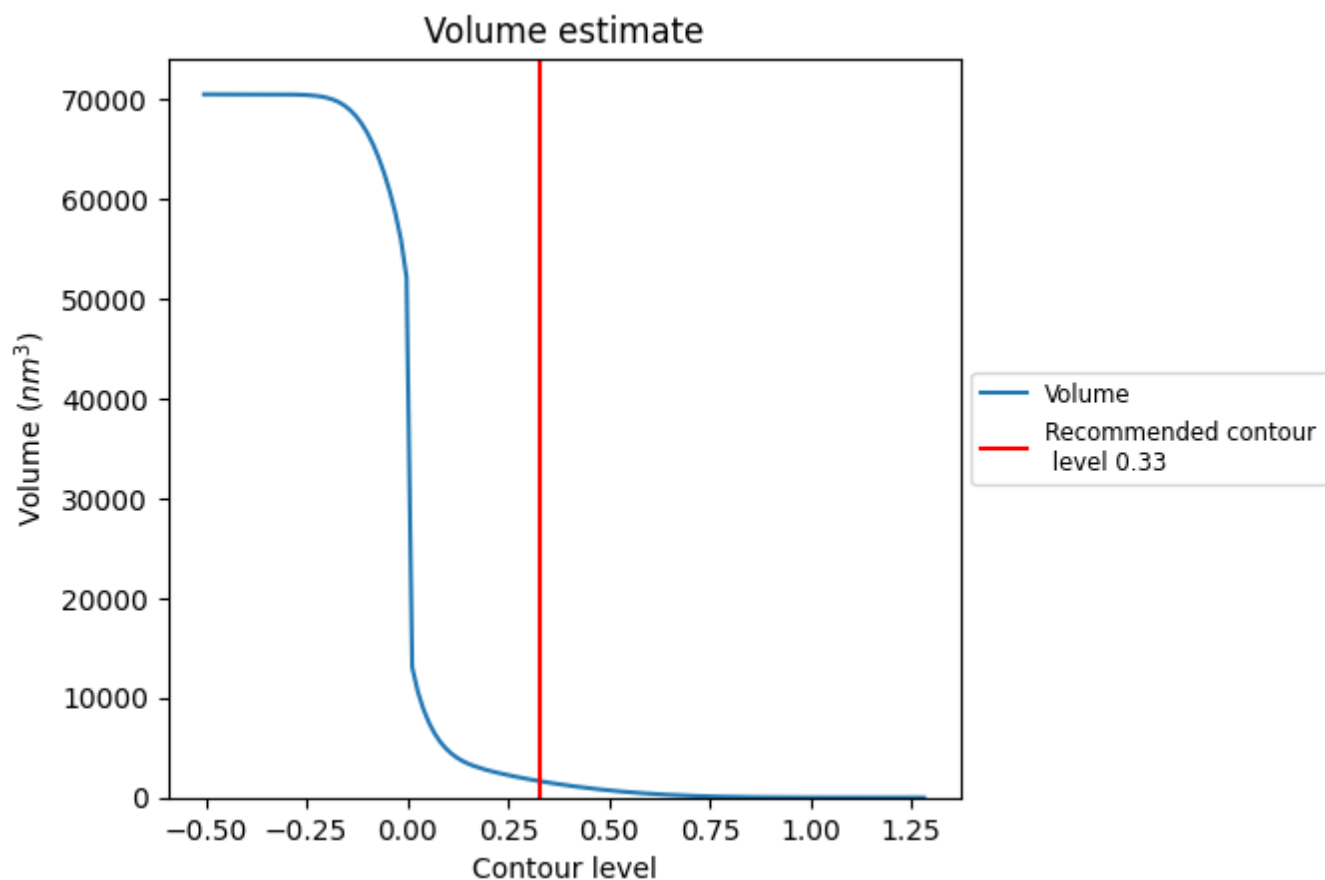
This section contains the results of statistical analysis of the map.

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.

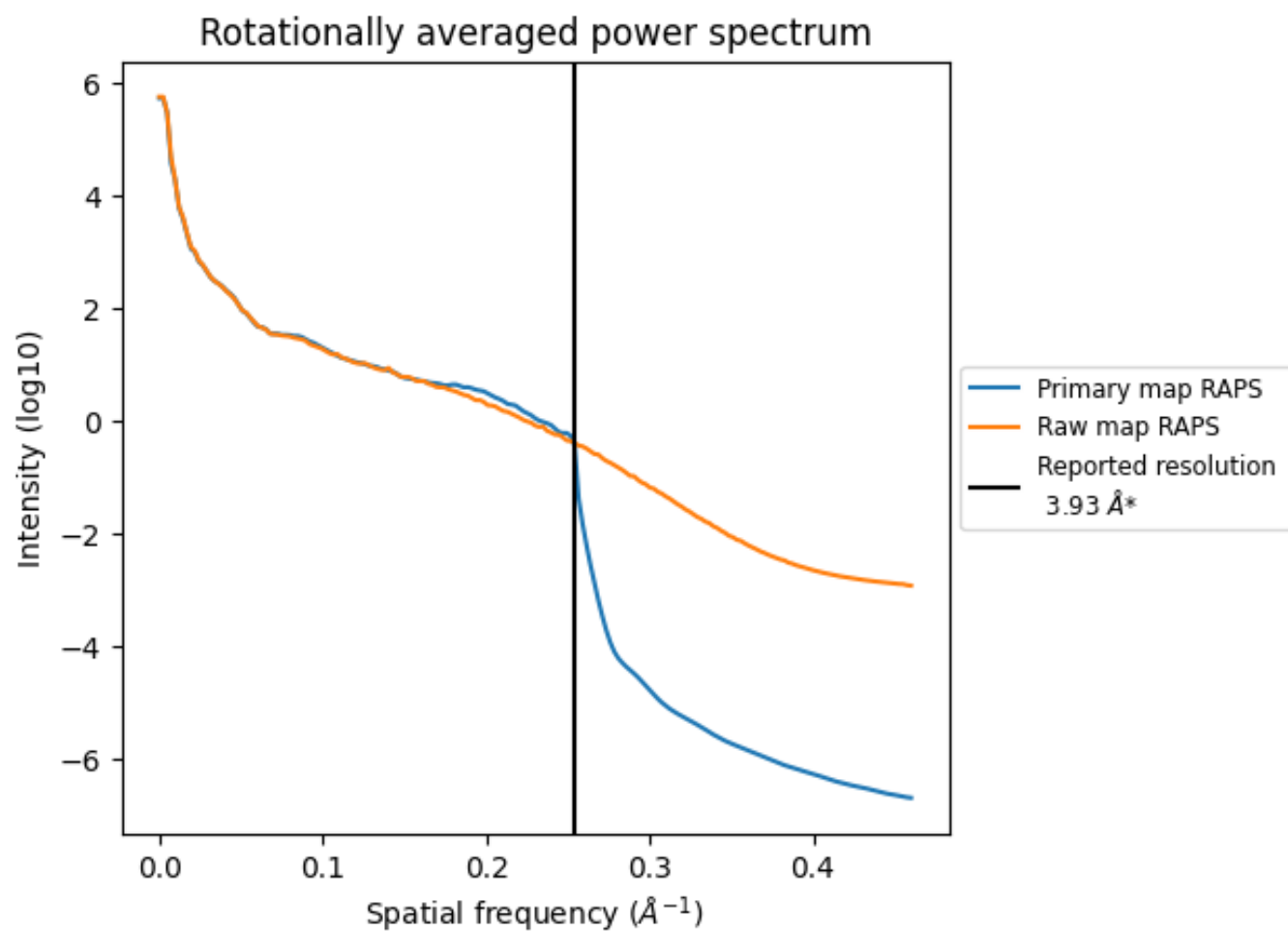
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1637 nm³; this corresponds to an approximate mass of 1479 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 ⓘ

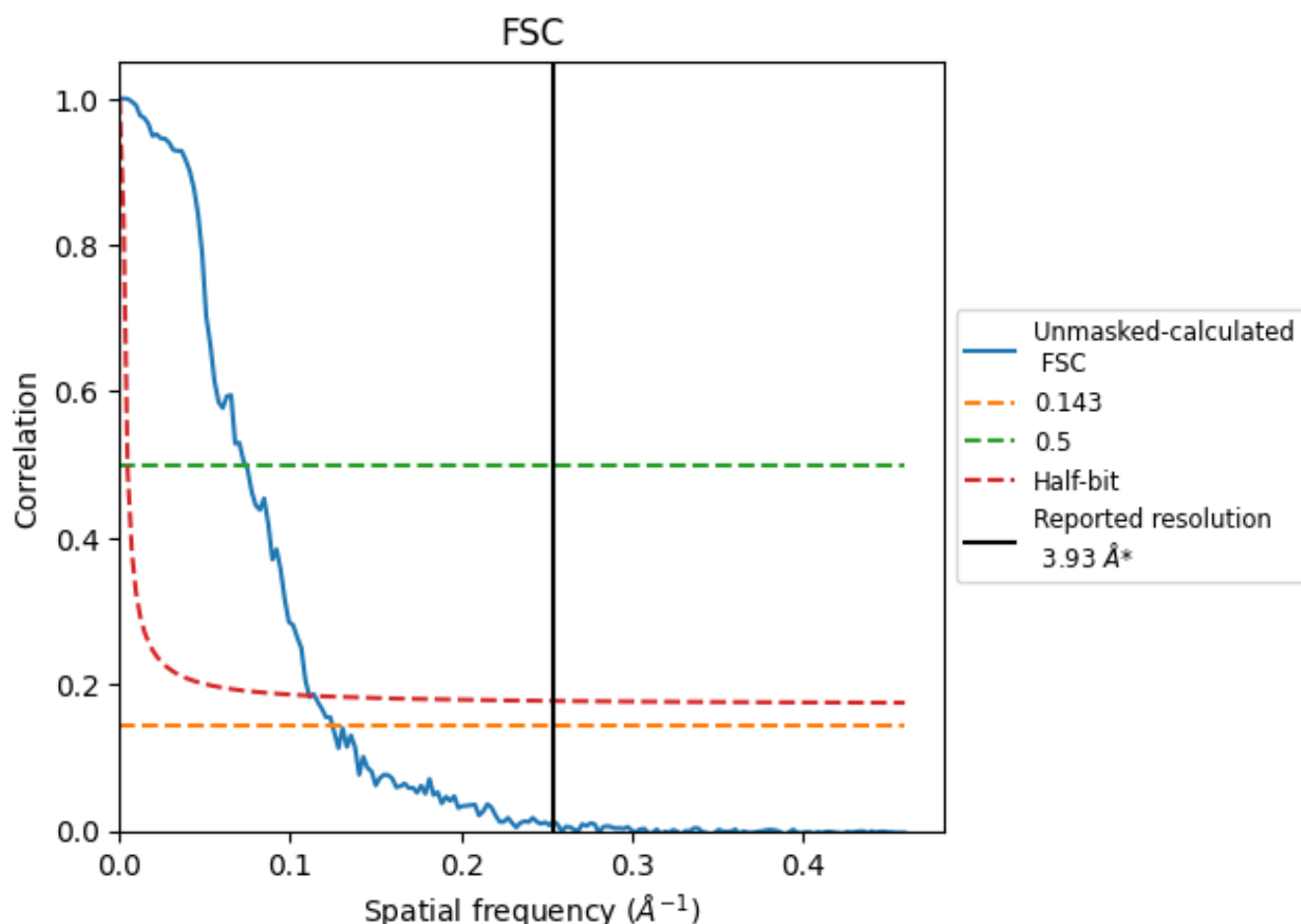


*Reported resolution corresponds to spatial frequency of 0.254 \AA^{-1}

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.254 \AA^{-1}

8.2 Resolution estimates [i](#)

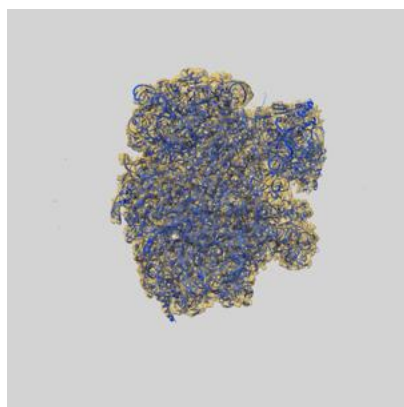
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.93	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	8.01	13.57	9.00

*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 8.01 differs from the reported value 3.93 by more than 10 %

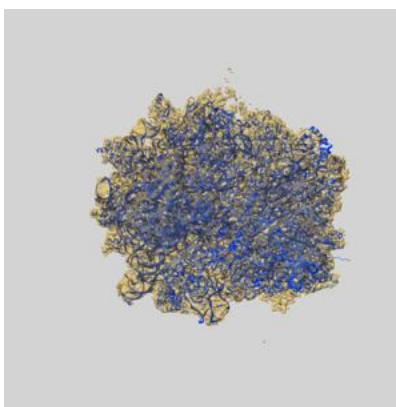
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-34866 and PDB model 8HL1. Per-residue inclusion information can be found in section [3](#) on page [17](#).

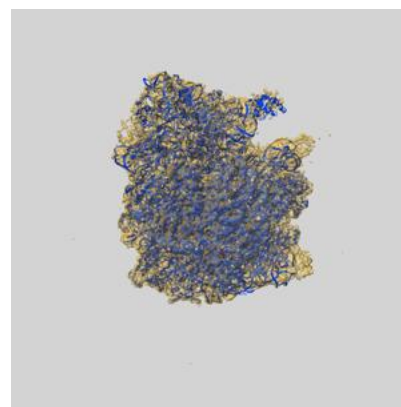
9.1 Map-model overlay [i](#)



X



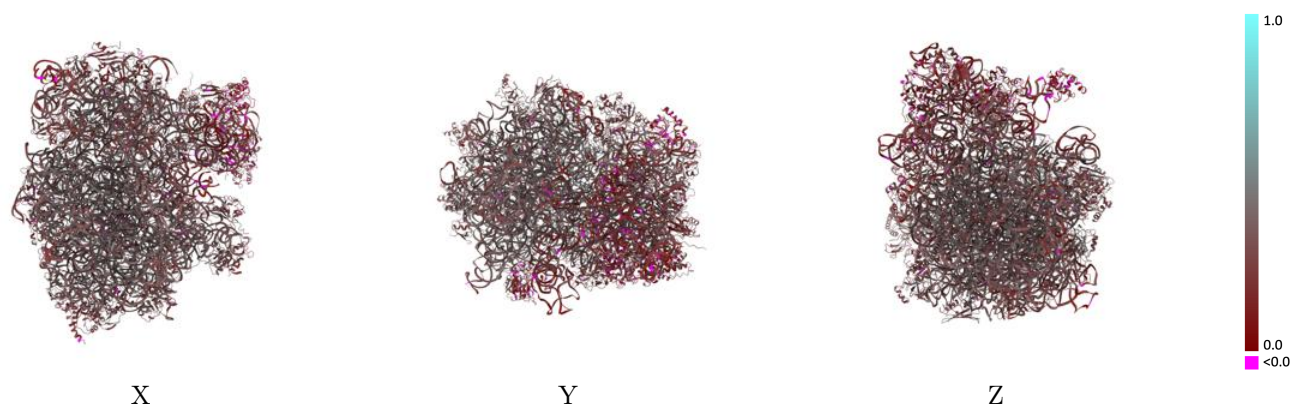
Y



Z

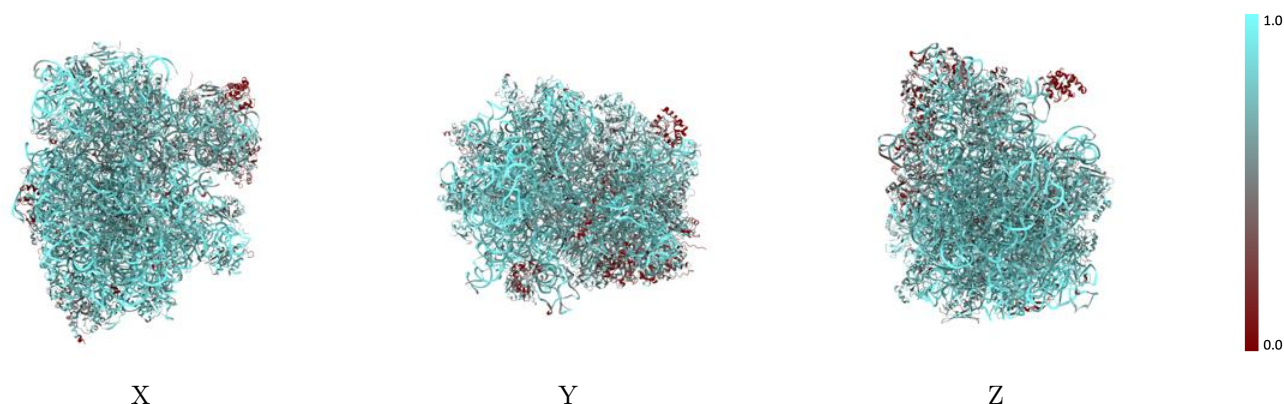
The images above show the 3D surface view of the map at the recommended contour level 0.33 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



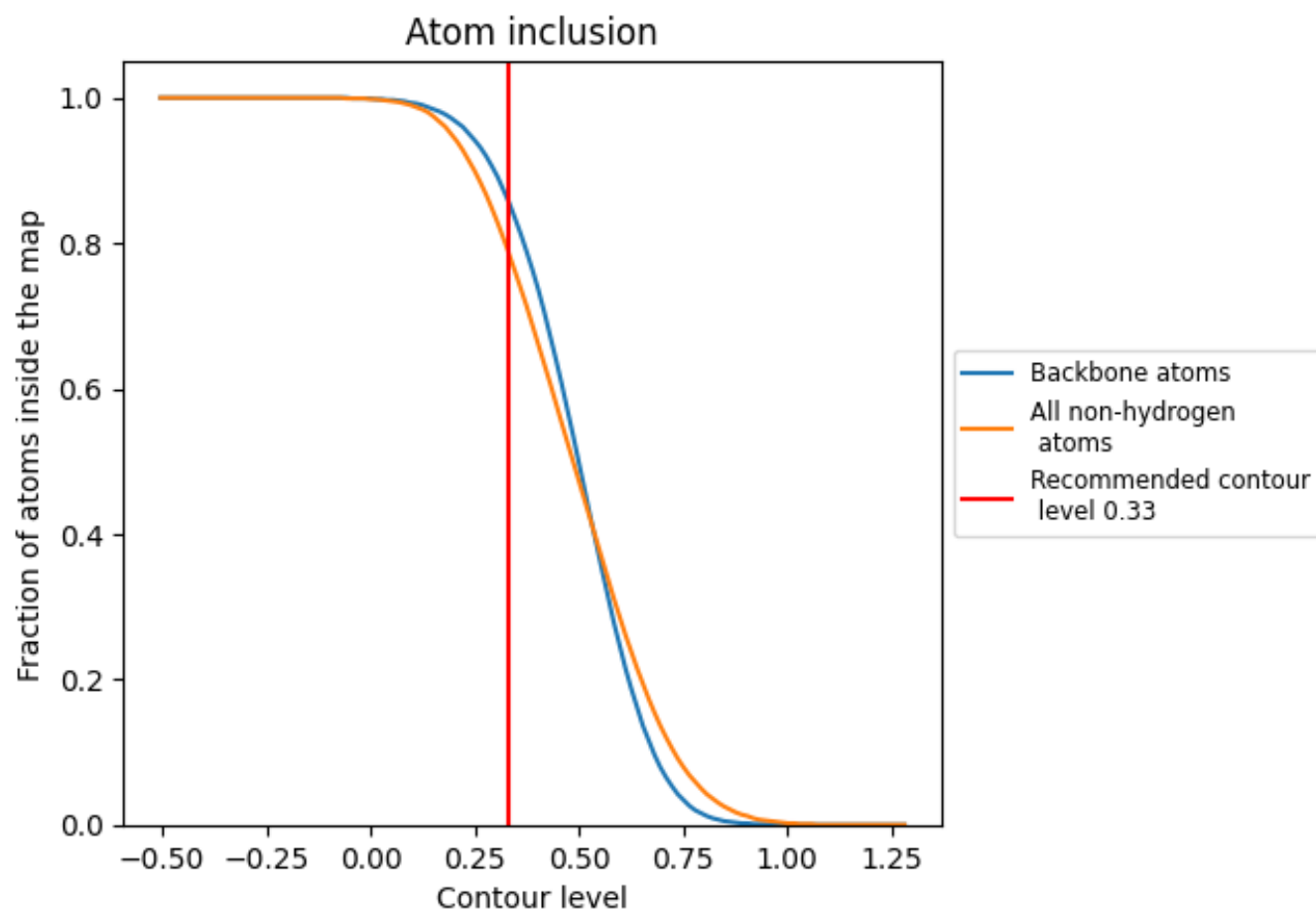
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.33).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 79% 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 (0.33) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7900	 0.3320
A16S	 0.8920	 0.3140
A23S	 0.9220	 0.3690
A5S	 0.9580	 0.3540
AEFG	 0.6340	 0.3100
AETN	 0.7700	 0.2280
AL1P	 0.3050	 0.1820
AL2P	 0.6650	 0.3980
AL3P	 0.7000	 0.3580
AL4P	 0.7020	 0.3600
AL5P	 0.6560	 0.2910
AL6P	 0.6880	 0.3430
ALX0	 0.6520	 0.3080
AMRN	 0.6500	 0.3000
APTN	 0.8690	 0.2930
APTP	 0.2690	 0.2070
AS2P	 0.4240	 0.2750
AS3P	 0.5340	 0.2690
AS4E	 0.6870	 0.3080
AS4P	 0.7090	 0.3020
AS5P	 0.6830	 0.3450
AS6E	 0.5920	 0.2800
AS7P	 0.3450	 0.1700
AS8E	 0.6290	 0.3220
AS8P	 0.7330	 0.3480
AS9P	 0.4420	 0.2080
L10E	 0.6620	 0.3610
L13P	 0.6690	 0.3260
L141	 0.6830	 0.3200
L142	 0.6670	 0.3130
L14P	 0.6160	 0.3870
L15E	 0.7020	 0.3830
L15P	 0.7420	 0.3790
L18E	 0.7540	 0.3600
L18P	 0.7260	 0.3190



Continued on next page...

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Chain	Atom inclusion	Q-score
L19E	 0.7170	 0.3670
L21E	 0.7050	 0.3960
L22P	 0.7150	 0.3680
L23P	 0.6410	 0.3620
L24E	 0.7720	 0.3800
L24P	 0.7120	 0.3470
L29P	 0.6330	 0.2920
L30E	 0.7190	 0.3320
L30P	 0.6880	 0.3130
L31E	 0.7180	 0.3670
L32E	 0.6250	 0.3610
L34E	 0.5760	 0.2580
L37A	 0.7090	 0.3850
L37E	 0.6520	 0.3720
L39E	 0.6660	 0.3430
L40E	 0.5180	 0.2910
L44E	 0.7530	 0.3840
L45A	 0.6190	 0.3150
L46A	 0.5350	 0.3500
L47A	 0.0670	 0.2470
L7A1	 0.6940	 0.3160
L7A2	 0.5280	 0.2820
S10P	 0.4600	 0.2320
S11P	 0.6110	 0.3220
S12P	 0.6080	 0.3690
S13P	 0.4020	 0.2010
S14P	 0.5680	 0.2560
S15P	 0.6960	 0.3120
S17E	 0.4950	 0.2390
S17P	 0.6830	 0.3650
S19E	 0.3690	 0.1860
S19P	 0.4940	 0.1980
S24E	 0.7350	 0.2920
S27A	 0.5650	 0.1530
S27E	 0.7720	 0.3140
S28E	 0.3930	 0.2580
S3AE	 0.6440	 0.2990
SL7A	 0.1540	 0.1530