



wwPDB EM Validation Summary Report ⓘ

Nov 26, 2022 – 10:22 PM EST

PDB ID : 7RYF
EMDB ID : EMD-24738
Title : A. baumannii Ribosome-TP-6076 complex: P-site tRNA 70S
Authors : Morgan, C.E.; Yu, E.W.
Deposited on : 2021-08-25
Resolution : 2.65 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

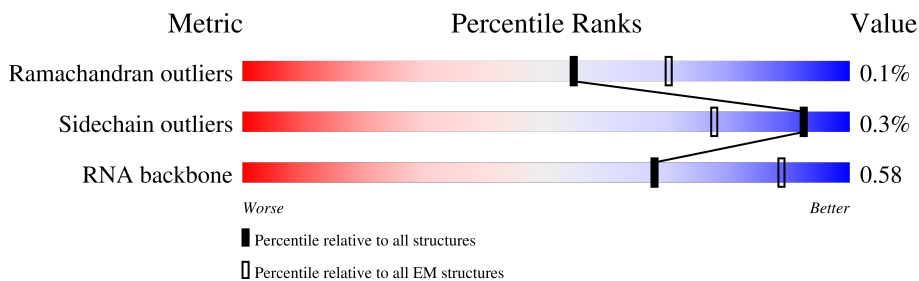
EMDB validation analysis : 0.0.1.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

1 Overall quality at a glance

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

The reported resolution of this entry is 2.65 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	0	51	
2	1	44	
3	2	64	
4	3	38	
5	A	2918	
6	B	115	
7	C	274	
8	D	212	

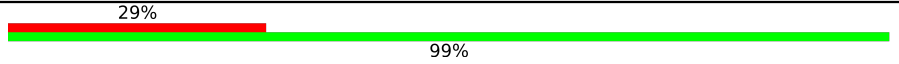
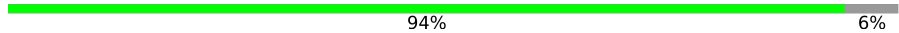


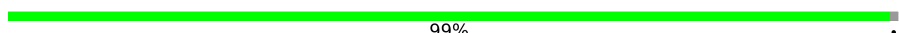
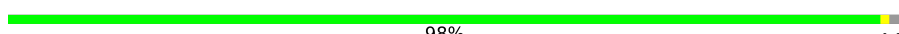
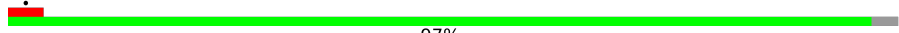







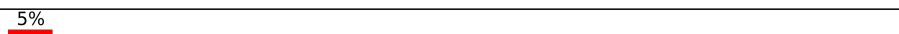
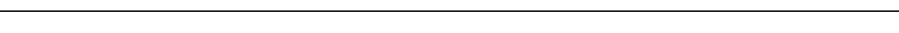
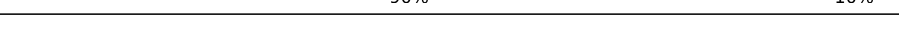
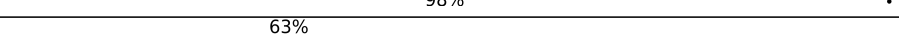


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Mol	Chain	Length	Quality of chain
9	E	200	6% 100%
10	F	178	95% 97% ..
11	G	177	72% 97% ..
12	H	148	26% 41% 59%
13	I	142	99% ..
14	J	122	100%
15	K	146	99% .
16	L	137	99% .
17	M	125	95% 5%
18	N	116	23% 98% .
19	O	122	7% 95% 5%
20	P	119	98% .
21	Q	103	100%
22	R	109	100%
23	S	106	9% 85% 15%
24	T	105	14% 98% .
25	U	98	7% 99% .
26	V	85	91% 9%
27	W	78	99% .
28	X	65	22% 92% 8%
29	Y	58	100%
30	Z	61	89% 11%
31	a	1544	11% 81% 17% ..
32	b	250	68% 89% 10%
33	c	250	84% 14%

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Mol	Chain	Length	Quality of chain
34	d	208	
35	e	165	
36	f	127	
37	g	156	
38	h	131	
39	i	128	
40	j	103	
41	k	128	
42	l	124	
43	m	118	
44	n	101	
45	o	89	
46	p	101	
47	q	85	
48	r	75	
49	s	91	
50	t	88	
51	u	71	
52	v	77	
53	w	3	

2 Entry composition [i](#)

There are 56 unique types of molecules in this entry. The entry contains 137516 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	0	48	400	256	72	69	3	0	0

- Molecule 2 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	1	44	363	222	85	54	2	0	0

- Molecule 3 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	2	63	509	319	110	76	4	0	0

- Molecule 4 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	3	38	295	179	64	48	4	0	0

- Molecule 5 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
5	A	2685	57610	25715	10555	18655	2685	0	0

- Molecule 6 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
6	B	115	2450	1095	440	800	115	0	0

- Molecule 7 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	C	271	2104	1297	435	364	8	0	0

- Molecule 8 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	D	210	1566	969	296	298	3	0	0

- Molecule 9 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	E	200	1516	952	281	278	5	0	0

- Molecule 10 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	F	174	1370	871	243	248	8	0	0

- Molecule 11 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	G	174	1318	832	236	249	1	0	0

- Molecule 12 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	H	60	458	287	84	86	1	0	0

- Molecule 13 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	I	141	1117	713	199	202	3	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
14	J	122	946	592	180	169	5	0	0

- Molecule 15 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
15	K	144	1071	663	213	195		0	0

- Molecule 16 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
16	L	137	1087	687	210	185	5	0	0

- Molecule 17 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
17	M	119	942	590	186	163	3	0	0

- Molecule 18 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
18	N	114	857	528	173	155	1	0	0

- Molecule 19 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
19	O	116	913	575	176	162		0	0

- Molecule 20 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
20	P	117	934	589	197	146	2	0	0

- Molecule 21 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	Q	103	Total	C	N	O	S	0	0
			807	506	155	143	3		

- Molecule 22 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	R	109	Total	C	N	O	S	0	0
			826	514	158	150	4		

- Molecule 23 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms				AltConf	Trace
23	S	90	Total	C	N	O	0	0
			702	447	127	128		

- Molecule 24 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms				AltConf	Trace
24	T	103	Total	C	N	O	0	0
			766	476	142	148		

- Molecule 25 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	U	97	Total	C	N	O	S	0	0
			760	477	143	139	1		

- Molecule 26 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	V	77	Total	C	N	O	S	0	0
			581	360	112	107	2		

- Molecule 27 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	W	77	Total	C	N	O	S	0	0
			632	395	130	105	2		

- Molecule 28 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	X	60	Total	C	N	O	S	0	0
			486	302	93	90	1		

- Molecule 29 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	Y	58	Total	C	N	O	S	0	0
			463	286	88	85	4		

- Molecule 30 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Z	54	Total	C	N	O	S	0	0
			447	265	100	81	1		

- Molecule 31 is a RNA chain called 16S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	a	1528	Total	C	N	O	P	0	0
			32782	14631	5994	10630	1527		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	1007	U	C	conflict	GB 1211343212
a	1034	C	U	conflict	GB 1211343212

- Molecule 32 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	b	225	Total	C	N	O	S	0	0
			1769	1110	328	325	6		

- Molecule 33 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	c	215	Total	C	N	O	S	0	0
			1690	1065	318	299	8		

- Molecule 34 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	d	207	Total	C	N	O	S	0	0
			1631	1017	313	299	2		

- Molecule 35 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	e	155	Total	C	N	O	S	0	0
			1129	700	217	207	5		

- Molecule 36 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	f	104	Total	C	N	O	S	0	0
			867	546	158	159	4		

- Molecule 37 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	g	136	Total	C	N	O	S	0	0
			1073	671	201	194	7		

- Molecule 38 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	h	130	Total	C	N	O	S	0	0
			985	615	177	187	6		

- Molecule 39 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	i	126	Total	C	N	O	S	0	0
			991	618	197	175	1		

- Molecule 40 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	j	100	Total	C	N	O	S	0	0
			801	500	150	148	3		

- Molecule 41 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	k	117	Total	C	N	O	S	0	0
			862	535	167	159	1		

- Molecule 42 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	l	122	Total	C	N	O	S	0	0
			945	580	193	167	5		

- Molecule 43 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	m	115	Total	C	N	O	S	0	0
			903	558	184	158	3		

- Molecule 44 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	n	100	Total	C	N	O	S	0	0
			792	493	158	137	4		

- Molecule 45 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	o	88	Total	C	N	O	S	0	0
			705	434	144	126	1		

- Molecule 46 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	p	82	Total	C	N	O	S	0	0
			644	403	128	112	1		

- Molecule 47 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	q	78	Total	C	N	O	S	0	0
			614	386	115	112	1		

- Molecule 48 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				AltConf	Trace
48	r	52	Total	C	N	O	0	0
			426	273	74	79		

- Molecule 49 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	s	82	Total	C	N	O	S	0	0
			646	412	125	107	2		

- Molecule 50 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	t	86	Total	C	N	O	S	0	0
			663	409	139	113	2		

- Molecule 51 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	u	57	Total	C	N	O	S	0	0
			473	294	96	82	1		

- Molecule 52 is a RNA chain called tRNA-met.

Mol	Chain	Residues	Atoms						AltConf	Trace
52	v	77	Total	C	N	O	P	S	0	0
			1636	733	291	535	76	1		

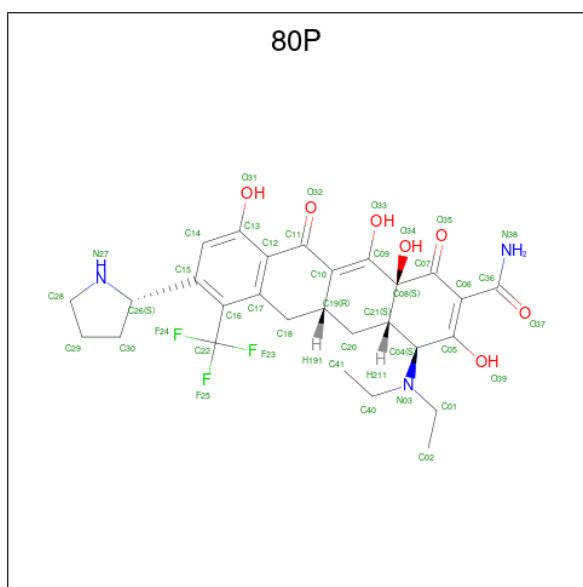
- Molecule 53 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	w	3	Total	C	N	O	P	0	0
			65	29	12	21	3		

- Molecule 54 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
54	3	1	Total	Zn	0
			1	1	

- Molecule 55 is (4S,4aS,5aR,12aS)-4-(diethylamino)-3,10,12,12a-tetrahydroxy-1,11-dioxo-8-[(2S)-pyrrolidin-2-yl]-7-(trifluoromethyl)-1,4,4a,5,5a,6,11,12a-octahydrotetracene-2-carboxamide (three-letter code: 80P) (formula: C₂₈H₃₂F₃N₃O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	F	N	O	
55	a	1	123	84	9	9	21	0
55	a	1	123	84	9	9	21	0
55	a	1	123	84	9	9	21	0

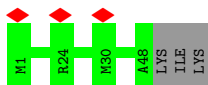
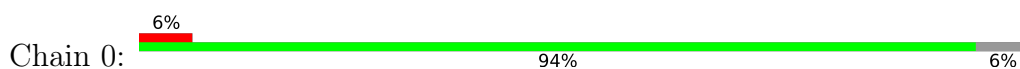
- Molecule 56 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
56	a	4	4	4	0

3 Residue-property plots [i](#)

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

- Molecule 1: 50S ribosomal protein L33

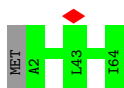


- Molecule 2: 50S ribosomal protein L34

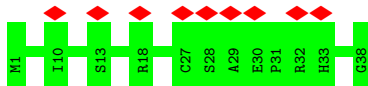


There are no outlier residues recorded for this chain.

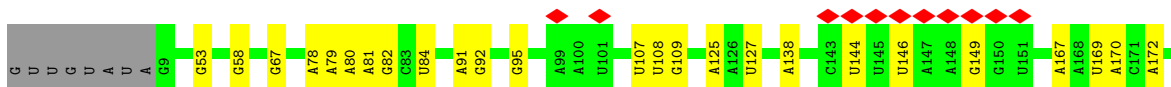
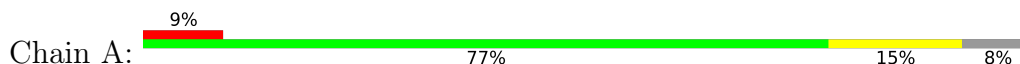
- Molecule 3: 50S ribosomal protein L35

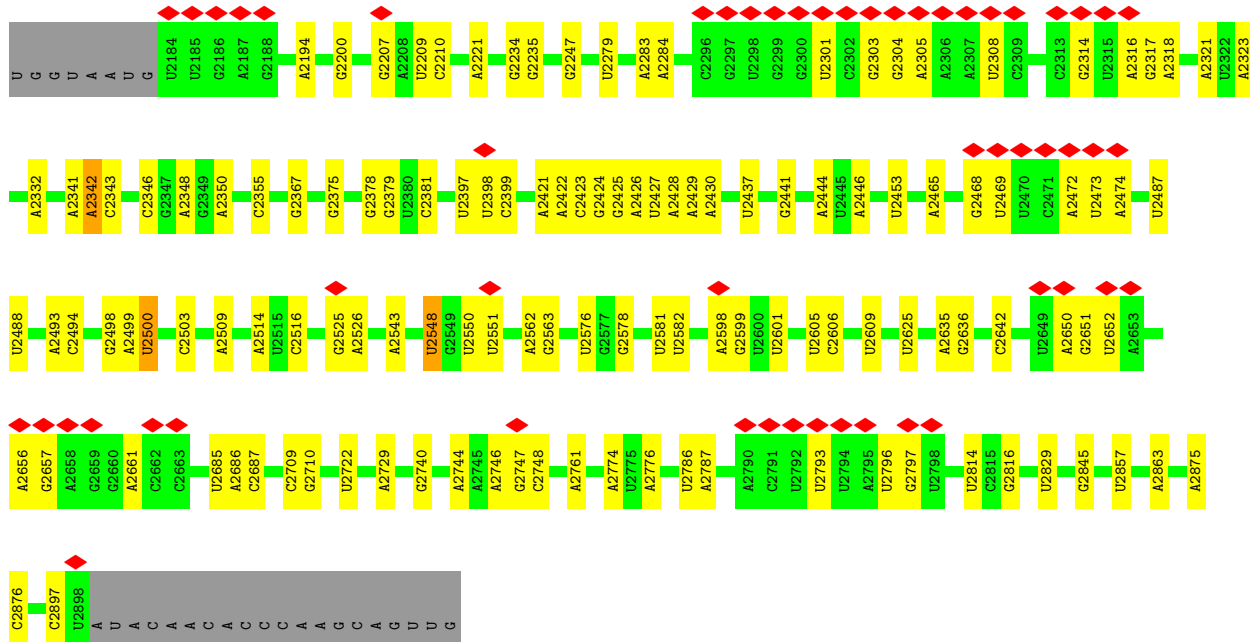


- Molecule 4: 50S ribosomal protein L36

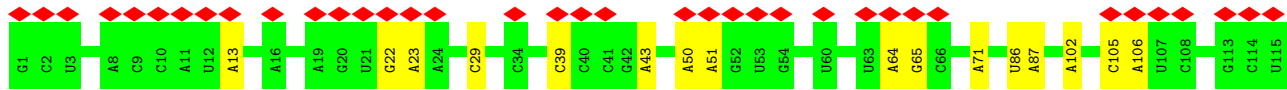
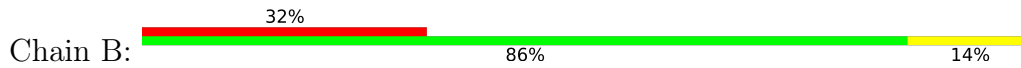


- Molecule 5: 23S ribosomal RNA

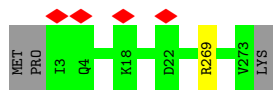




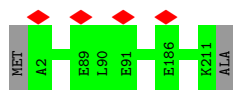
• Molecule 6: 5S ribosomal RNA



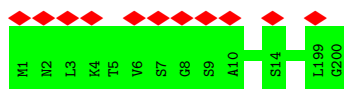
• Molecule 7: 50S ribosomal protein L2



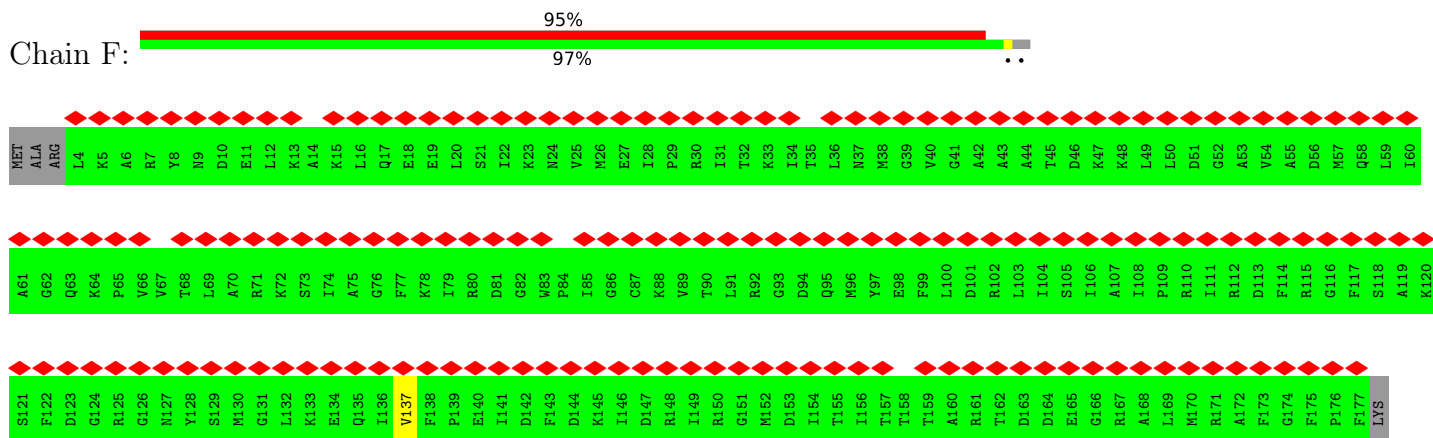
• Molecule 8: 50S ribosomal protein L3



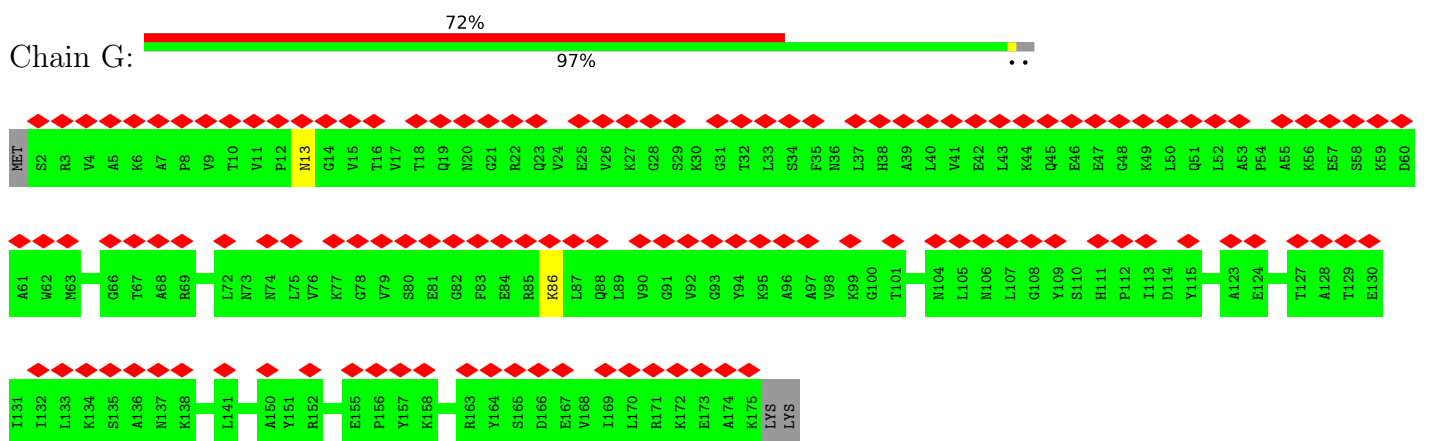
• Molecule 9: 50S ribosomal protein L4



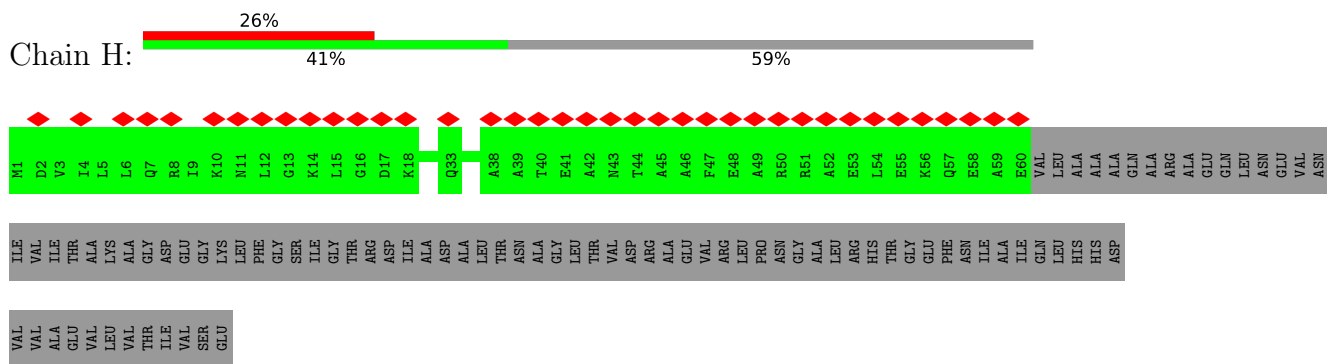
• Molecule 10: 50S ribosomal protein L5



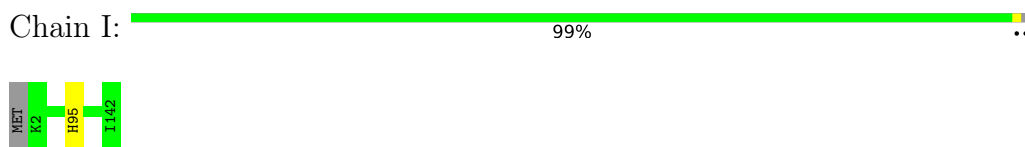
• Molecule 11: 50S ribosomal protein L6



• Molecule 12: 50S ribosomal protein L9

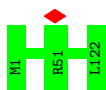


• Molecule 13: 50S ribosomal protein L13



• Molecule 14: 50S ribosomal protein L14





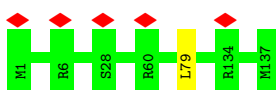
- Molecule 15: 50S ribosomal protein L15

Chain K: 99%



- Molecule 16: 50S ribosomal protein L16

Chain L: 99%



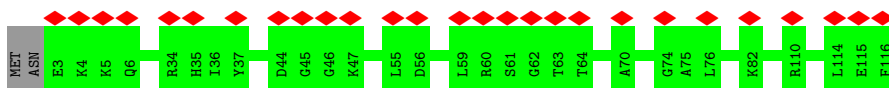
- Molecule 17: 50S ribosomal protein L17

Chain M: 95% 5%



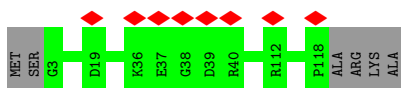
- Molecule 18: 50S ribosomal protein L18

Chain N: 23% 98%



- Molecule 19: 50S ribosomal protein L19

Chain O: 7% 95% 5%



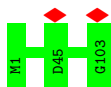
- Molecule 20: 50S ribosomal protein L20

Chain P: 98%



- Molecule 21: 50S ribosomal protein L21

Chain Q:  100%




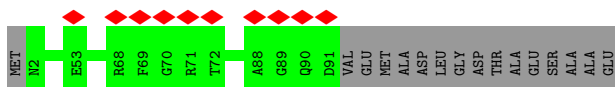
- Molecule 22: 50S ribosomal protein L22

Chain R:  100%



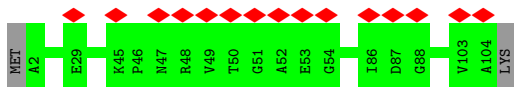
- Molecule 23: 50S ribosomal protein L23

Chain S:  9% 85% 15%



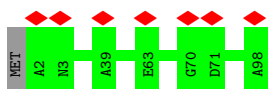
- Molecule 24: 50S ribosomal protein L24

Chain T:  14% 98%

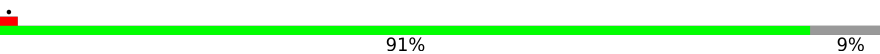


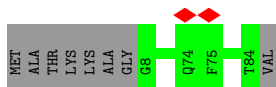
- Molecule 25: 50S ribosomal protein L25

Chain U:  7% 99%



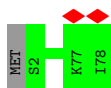
- Molecule 26: 50S ribosomal protein L27

Chain V:  91% 9%

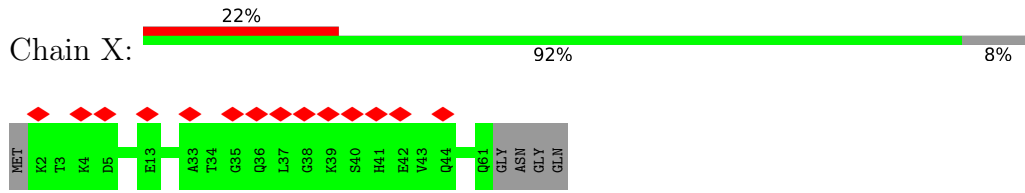


- Molecule 27: 50S ribosomal protein L28

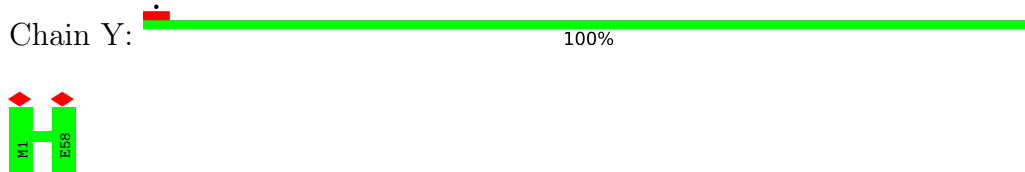
Chain W:  99%



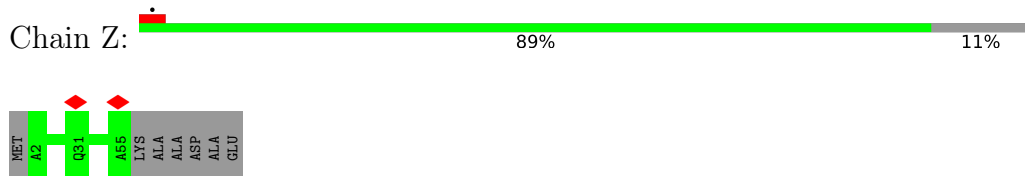
- Molecule 28: 50S ribosomal protein L29



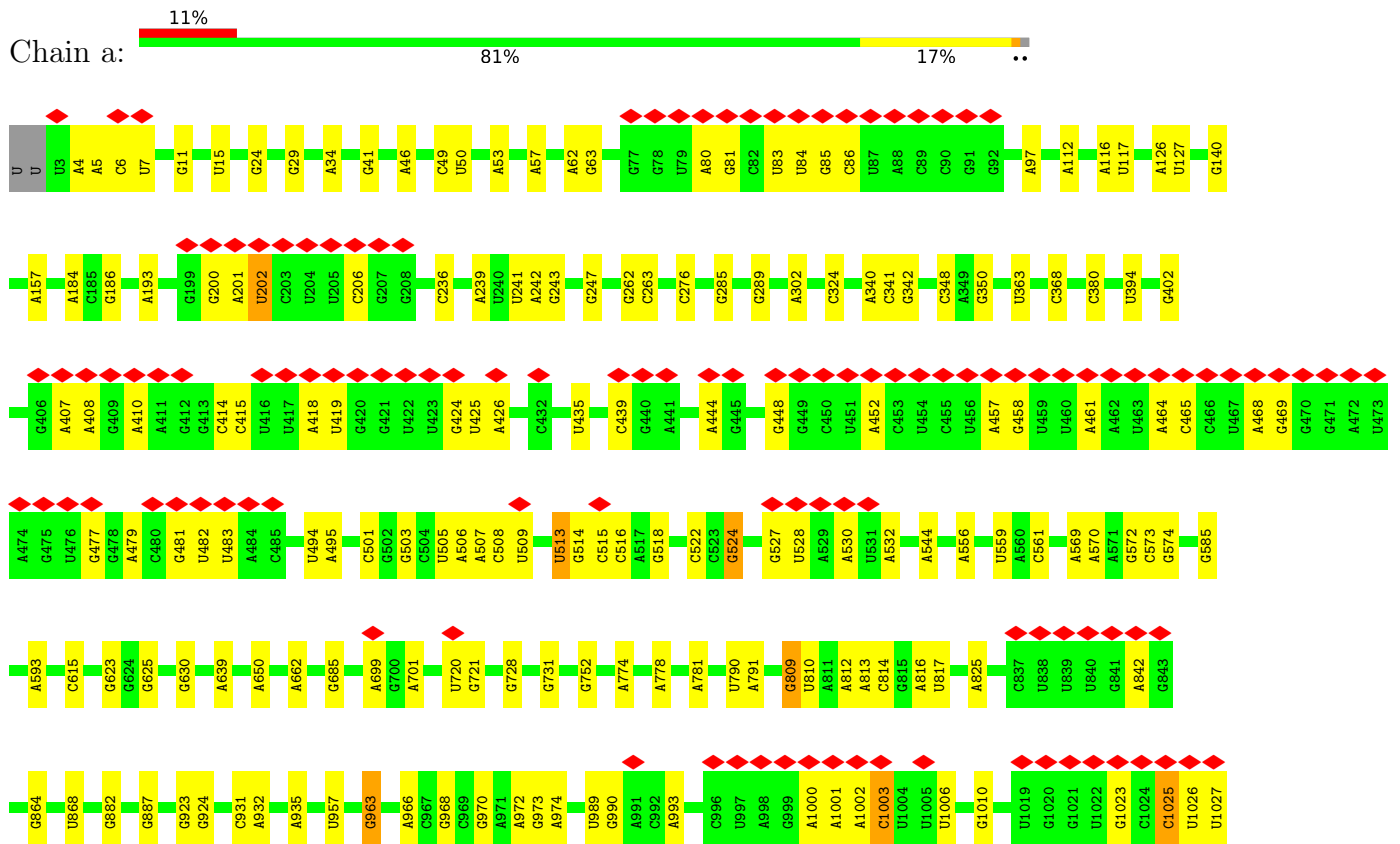
- Molecule 29: 50S ribosomal protein L30

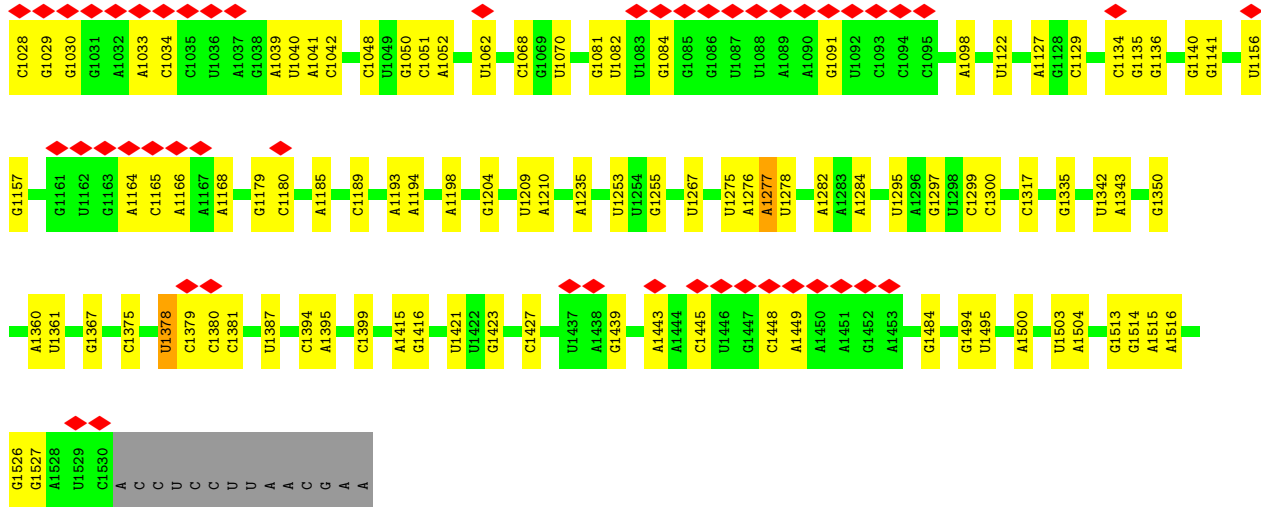


- Molecule 30: 50S ribosomal protein L32

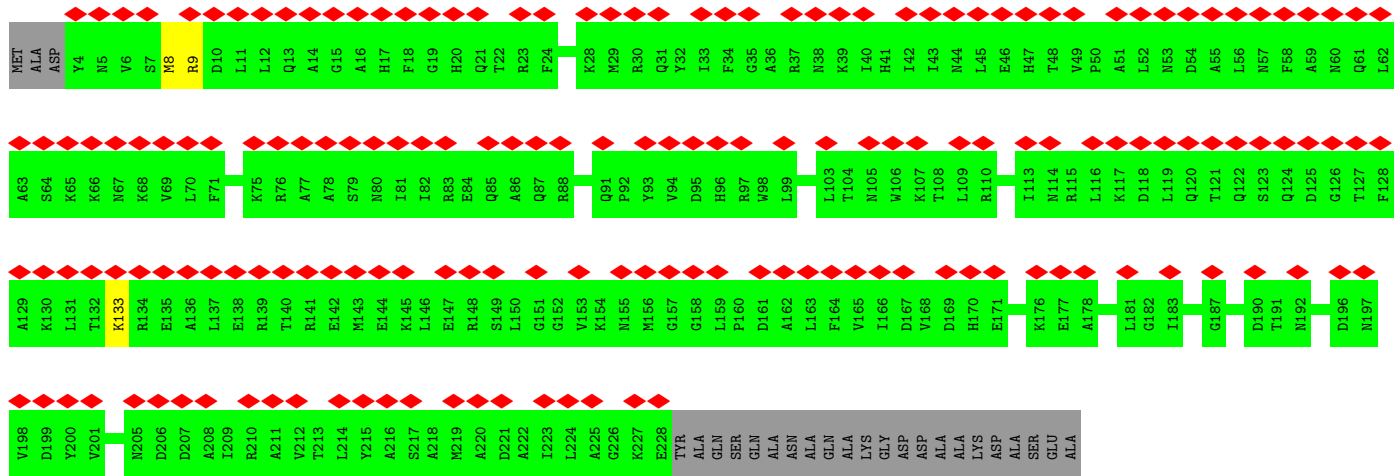
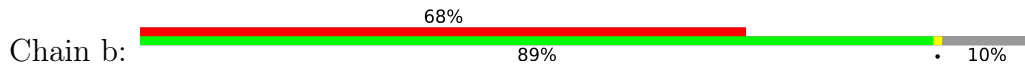


- Molecule 31: 16S Ribosomal RNA

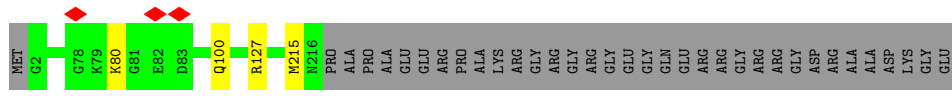
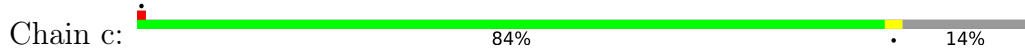




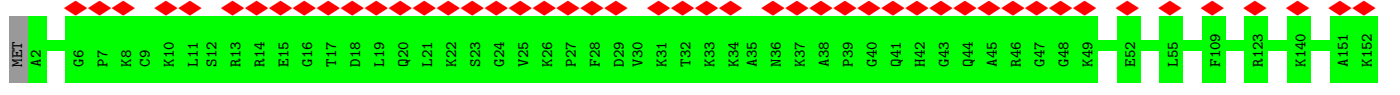
• Molecule 32: 30S ribosomal protein S2

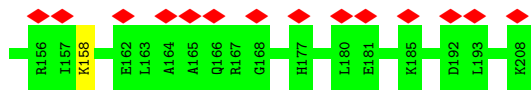


• Molecule 33: 30S ribosomal protein S3

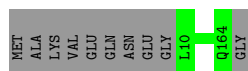


• Molecule 34: 30S ribosomal protein S4

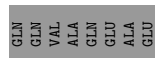
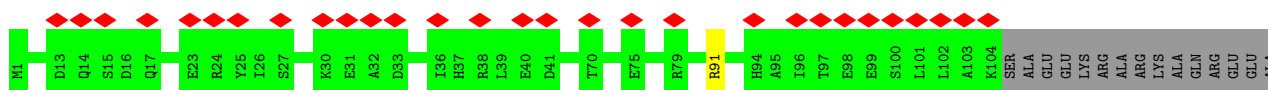
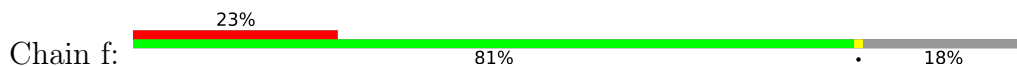




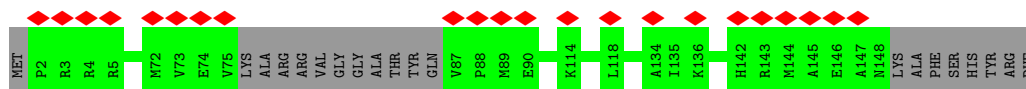
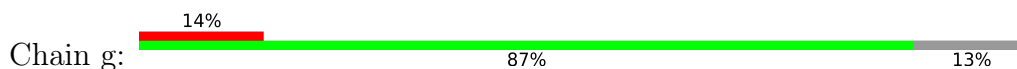
• Molecule 35: 30S ribosomal protein S5



• Molecule 36: 30S ribosomal protein S6



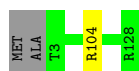
• Molecule 37: 30S ribosomal protein S7



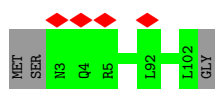
• Molecule 38: 30S ribosomal protein S8



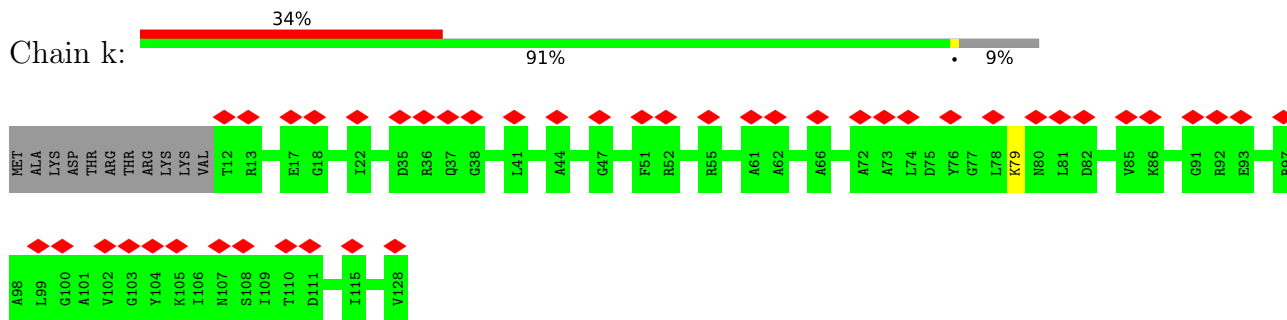
• Molecule 39: 30S ribosomal protein S9



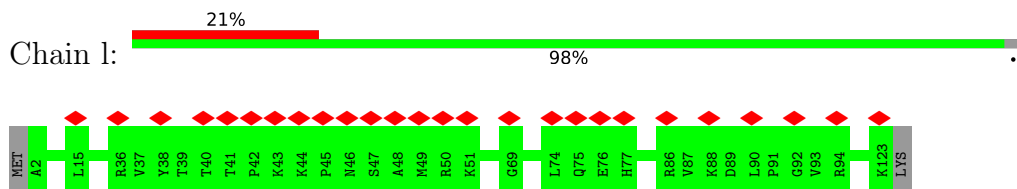
• Molecule 40: 30S ribosomal protein S10



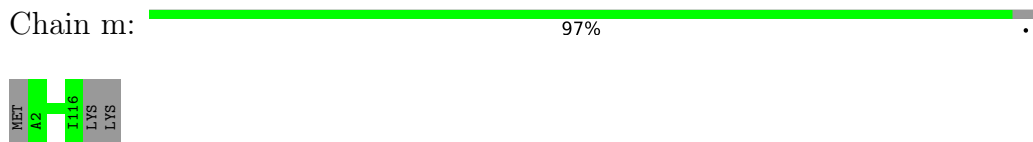
• Molecule 41: 30S ribosomal protein S11



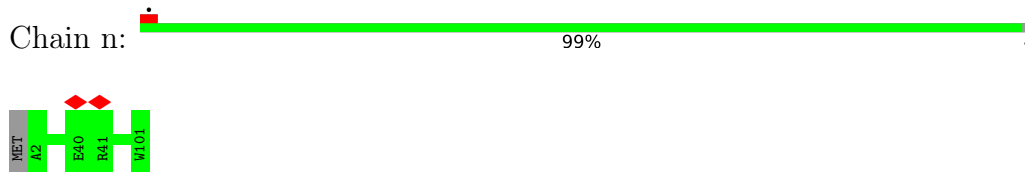
• Molecule 42: 30S ribosomal protein S12



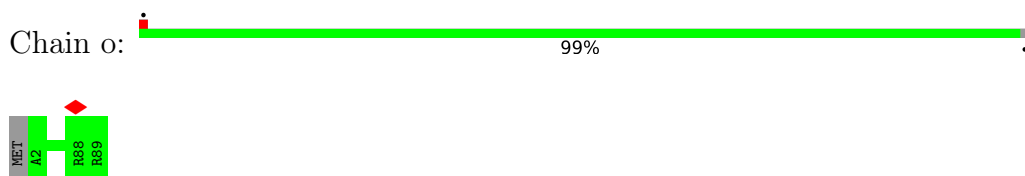
• Molecule 43: 30S ribosomal protein S13



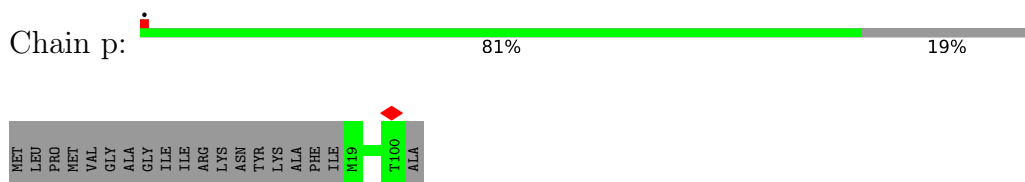
• Molecule 44: 30S ribosomal protein S14



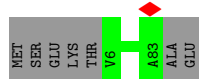
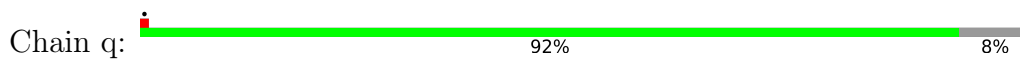
• Molecule 45: 30S ribosomal protein S15



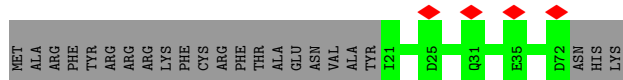
• Molecule 46: 30S ribosomal protein S16



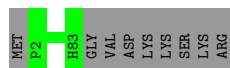
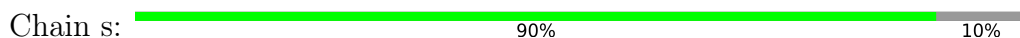
• Molecule 47: 30S ribosomal protein S17



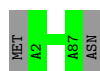
• Molecule 48: 30S ribosomal protein S18



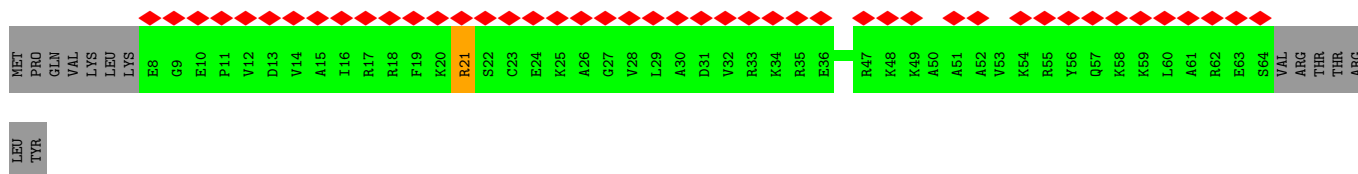
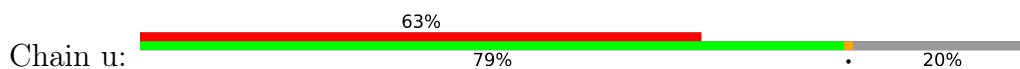
• Molecule 49: 30S ribosomal protein S19



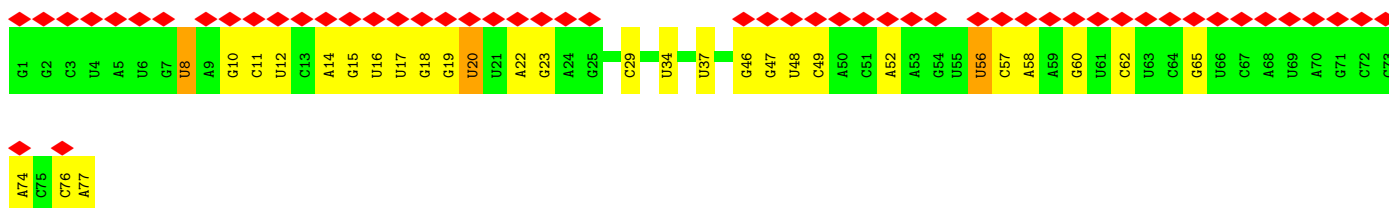
• Molecule 50: 30S ribosomal protein S20



• Molecule 51: 30S ribosomal protein S21



• Molecule 52: tRNA-met



• Molecule 53: mRNA

Chain w:  100%

There are no outlier residues recorded for this chain.

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	47011	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	4.322	Depositor
Minimum map value	-0.094	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.039	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	557.568, 557.568, 557.568	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.089, 1.089, 1.089	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: H2U, UR3, 5MU, 80P, 6MZ, 2MG, 7MG, MG, ZN, 2MA, MA6, OMG, OMU, 5MC, 4OC, 4SU, PSU

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	0	0.27	0/407	0.51	0/540
2	1	0.24	0/367	0.66	0/481
3	2	0.24	0/515	0.61	0/678
4	3	0.25	0/296	0.59	0/389
5	A	0.21	0/64227	0.76	3/100172 (0.0%)
6	B	0.18	0/2739	0.74	0/4266
7	C	0.24	0/2144	0.58	0/2880
8	D	0.24	0/1584	0.56	0/2135
9	E	0.25	0/1537	0.51	0/2073
10	F	0.26	0/1390	0.59	0/1863
11	G	0.27	0/1337	0.53	0/1807
12	H	0.27	0/461	0.53	0/616
13	I	0.25	0/1143	0.49	0/1541
14	J	0.25	0/956	0.59	0/1286
15	K	0.24	0/1079	0.58	0/1439
16	L	0.25	0/1104	0.58	0/1475
17	M	0.24	0/956	0.57	0/1282
18	N	0.26	0/865	0.58	0/1156
19	O	0.25	0/925	0.57	0/1241
20	P	0.26	0/947	0.59	0/1262
21	Q	0.24	0/818	0.55	0/1094
22	R	0.24	0/831	0.51	0/1113
23	S	0.25	0/708	0.54	0/947
24	T	0.24	0/770	0.51	0/1034
25	U	0.25	0/770	0.55	0/1036
26	V	0.25	0/589	0.53	0/788
27	W	0.24	0/642	0.56	0/856
28	X	0.24	0/487	0.47	0/646
29	Y	0.26	0/468	0.53	0/624
30	Z	0.24	0/453	0.63	0/604
31	a	0.29	0/36476	0.77	10/56895 (0.0%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	b	0.25	0/1799	0.52	0/2429
33	c	0.28	0/1714	0.59	1/2304 (0.0%)
34	d	0.27	0/1653	0.56	0/2213
35	e	0.26	0/1141	0.58	0/1537
36	f	0.25	0/882	0.53	0/1189
37	g	0.26	0/1087	0.56	0/1456
38	h	0.25	0/993	0.52	0/1331
39	i	0.27	0/1002	0.59	0/1339
40	j	0.26	0/811	0.54	0/1096
41	k	0.24	0/878	0.52	0/1189
42	l	0.24	0/958	0.58	0/1284
43	m	0.25	0/913	0.59	0/1226
44	n	0.29	0/803	0.61	0/1071
45	o	0.25	0/715	0.52	0/958
46	p	0.25	0/655	0.56	0/879
47	q	0.24	0/621	0.53	0/837
48	r	0.25	0/432	0.51	0/583
49	s	0.27	0/664	0.52	0/897
50	t	0.26	0/669	0.52	0/892
51	u	0.30	0/478	0.63	1/631 (0.2%)
52	v	0.19	0/1739	0.76	0/2709
53	w	0.15	0/72	0.72	0/110
All	All	0.25	0/148670	0.71	15/222379 (0.0%)

There are no bond length outliers.

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	a	415	C	N3-C2-O2	-7.00	117.00	121.90
31	a	1379	C	C2-N1-C1'	6.81	126.29	118.80
31	a	202	U	C2-N1-C1'	6.47	125.47	117.70
31	a	1378	U	C2-N1-C1'	6.18	125.12	117.70
5	A	1583	C	C2-N1-C1'	5.80	125.18	118.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	0	46/51 (90%)	45 (98%)	1 (2%)	0	100	100
2	1	42/44 (96%)	40 (95%)	2 (5%)	0	100	100
3	2	61/64 (95%)	59 (97%)	2 (3%)	0	100	100
4	3	36/38 (95%)	36 (100%)	0	0	100	100
7	C	269/274 (98%)	261 (97%)	8 (3%)	0	100	100
8	D	208/212 (98%)	200 (96%)	8 (4%)	0	100	100
9	E	198/200 (99%)	197 (100%)	1 (0%)	0	100	100
10	F	172/178 (97%)	163 (95%)	8 (5%)	1 (1%)	25	37
11	G	172/177 (97%)	170 (99%)	1 (1%)	1 (1%)	25	37
12	H	58/148 (39%)	56 (97%)	2 (3%)	0	100	100
13	I	139/142 (98%)	138 (99%)	1 (1%)	0	100	100
14	J	120/122 (98%)	118 (98%)	2 (2%)	0	100	100
15	K	142/146 (97%)	137 (96%)	5 (4%)	0	100	100
16	L	135/137 (98%)	132 (98%)	2 (2%)	1 (1%)	22	33
17	M	117/125 (94%)	115 (98%)	2 (2%)	0	100	100
18	N	112/116 (97%)	112 (100%)	0	0	100	100
19	O	114/122 (93%)	112 (98%)	2 (2%)	0	100	100
20	P	115/119 (97%)	115 (100%)	0	0	100	100
21	Q	101/103 (98%)	97 (96%)	4 (4%)	0	100	100
22	R	107/109 (98%)	103 (96%)	4 (4%)	0	100	100
23	S	88/106 (83%)	88 (100%)	0	0	100	100
24	T	101/105 (96%)	96 (95%)	5 (5%)	0	100	100
25	U	95/98 (97%)	93 (98%)	2 (2%)	0	100	100
26	V	75/85 (88%)	74 (99%)	1 (1%)	0	100	100
27	W	75/78 (96%)	75 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
28	X	58/65 (89%)	58 (100%)	0	0	100	100
29	Y	56/58 (97%)	55 (98%)	1 (2%)	0	100	100
30	Z	52/61 (85%)	51 (98%)	1 (2%)	0	100	100
32	b	223/250 (89%)	211 (95%)	11 (5%)	1 (0%)	34	48
33	c	213/250 (85%)	202 (95%)	11 (5%)	0	100	100
34	d	205/208 (99%)	198 (97%)	7 (3%)	0	100	100
35	e	153/165 (93%)	151 (99%)	2 (1%)	0	100	100
36	f	102/127 (80%)	96 (94%)	6 (6%)	0	100	100
37	g	132/156 (85%)	130 (98%)	2 (2%)	0	100	100
38	h	128/131 (98%)	124 (97%)	4 (3%)	0	100	100
39	i	124/128 (97%)	121 (98%)	3 (2%)	0	100	100
40	j	98/103 (95%)	94 (96%)	4 (4%)	0	100	100
41	k	115/128 (90%)	112 (97%)	3 (3%)	0	100	100
42	l	120/124 (97%)	115 (96%)	5 (4%)	0	100	100
43	m	113/118 (96%)	109 (96%)	4 (4%)	0	100	100
44	n	98/101 (97%)	94 (96%)	4 (4%)	0	100	100
45	o	86/89 (97%)	85 (99%)	1 (1%)	0	100	100
46	p	80/101 (79%)	80 (100%)	0	0	100	100
47	q	76/85 (89%)	76 (100%)	0	0	100	100
48	r	50/75 (67%)	50 (100%)	0	0	100	100
49	s	80/91 (88%)	80 (100%)	0	0	100	100
50	t	84/88 (96%)	84 (100%)	0	0	100	100
51	u	55/71 (78%)	52 (94%)	3 (6%)	0	100	100
All	All	5399/5872 (92%)	5260 (97%)	135 (2%)	4 (0%)	54	69

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	F	137	VAL
11	G	13	ASN
16	L	79	LEU
32	b	8	MET

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	0	44/47 (94%)	44 (100%)	0	100	100
2	1	36/36 (100%)	36 (100%)	0	100	100
3	2	52/53 (98%)	52 (100%)	0	100	100
4	3	33/33 (100%)	33 (100%)	0	100	100
7	C	217/220 (99%)	216 (100%)	1 (0%)	88	94
8	D	166/167 (99%)	166 (100%)	0	100	100
9	E	155/155 (100%)	155 (100%)	0	100	100
10	F	144/147 (98%)	144 (100%)	0	100	100
11	G	139/142 (98%)	138 (99%)	1 (1%)	84	91
12	H	45/112 (40%)	45 (100%)	0	100	100
13	I	117/118 (99%)	116 (99%)	1 (1%)	78	87
14	J	103/103 (100%)	103 (100%)	0	100	100
15	K	106/108 (98%)	106 (100%)	0	100	100
16	L	113/113 (100%)	113 (100%)	0	100	100
17	M	96/101 (95%)	96 (100%)	0	100	100
18	N	83/85 (98%)	83 (100%)	0	100	100
19	O	98/102 (96%)	98 (100%)	0	100	100
20	P	85/86 (99%)	85 (100%)	0	100	100
21	Q	84/84 (100%)	84 (100%)	0	100	100
22	R	88/88 (100%)	88 (100%)	0	100	100
23	S	76/87 (87%)	76 (100%)	0	100	100
24	T	83/85 (98%)	83 (100%)	0	100	100
25	U	79/80 (99%)	79 (100%)	0	100	100
26	V	59/64 (92%)	59 (100%)	0	100	100
27	W	69/70 (99%)	69 (100%)	0	100	100
28	X	53/56 (95%)	53 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
29	Y	54/54 (100%)	54 (100%)	0	100	100
30	Z	46/50 (92%)	46 (100%)	0	100	100
32	b	185/200 (92%)	183 (99%)	2 (1%)	73	85
33	c	175/198 (88%)	172 (98%)	3 (2%)	60	77
34	d	170/171 (99%)	169 (99%)	1 (1%)	86	92
35	e	113/120 (94%)	113 (100%)	0	100	100
36	f	94/111 (85%)	93 (99%)	1 (1%)	73	85
37	g	113/128 (88%)	113 (100%)	0	100	100
38	h	108/109 (99%)	108 (100%)	0	100	100
39	i	99/100 (99%)	98 (99%)	1 (1%)	76	86
40	j	89/91 (98%)	89 (100%)	0	100	100
41	k	88/98 (90%)	87 (99%)	1 (1%)	73	85
42	l	104/106 (98%)	104 (100%)	0	100	100
43	m	95/98 (97%)	95 (100%)	0	100	100
44	n	81/82 (99%)	81 (100%)	0	100	100
45	o	71/72 (99%)	71 (100%)	0	100	100
46	p	63/77 (82%)	63 (100%)	0	100	100
47	q	70/76 (92%)	70 (100%)	0	100	100
48	r	46/66 (70%)	46 (100%)	0	100	100
49	s	70/78 (90%)	70 (100%)	0	100	100
50	t	65/67 (97%)	65 (100%)	0	100	100
51	u	48/62 (77%)	47 (98%)	1 (2%)	53	72
All	All	4470/4756 (94%)	4457 (100%)	13 (0%)	92	96

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
33	c	127	ARG
34	d	158	LYS
51	u	21	ARG
39	i	104	ARG
41	k	79	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
32	b	61	GLN
37	g	29	HIS
40	j	56	HIS
16	L	46	GLN
9	E	162	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
31	a	1524/1544 (98%)	267 (17%)	0
5	A	2677/2918 (91%)	412 (15%)	37 (1%)
52	v	76/77 (98%)	30 (39%)	0
53	w	2/3 (66%)	0	0
6	B	114/115 (99%)	15 (13%)	2 (1%)
All	All	4393/4657 (94%)	724 (16%)	39 (0%)

5 of 724 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	A	53	G
5	A	58	G
5	A	67	G
5	A	78	A
5	A	80	A

5 of 39 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	A	1816	U
5	A	2651	G
5	A	1880	G
5	A	2342	A
6	B	50	A

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

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
52	PSU	v	56	52	18,21,22	1.09	1 (5%)	22,30,33	1.79	5 (22%)
5	PSU	A	2576	5	18,21,22	1.12	1 (5%)	22,30,33	1.77	5 (22%)
52	5MU	v	55	52	19,22,23	0.40	0	28,32,35	0.57	0
31	2MG	a	963	31	18,26,27	1.14	2 (11%)	16,38,41	1.00	1 (6%)
31	PSU	a	513	31	18,21,22	1.05	1 (5%)	22,30,33	1.72	3 (13%)
5	PSU	A	2500	5	18,21,22	1.11	1 (5%)	22,30,33	1.81	5 (22%)
31	4OC	a	1399	31	20,23,24	3.16	8 (40%)	26,32,35	0.88	1 (3%)
5	6MZ	A	2026	5	18,25,26	1.84	3 (16%)	16,36,39	3.39	5 (31%)
5	OMG	A	2247	52,5	18,26,27	1.11	2 (11%)	19,38,41	0.86	1 (5%)
5	PSU	A	2601	5	18,21,22	1.06	1 (5%)	22,30,33	1.78	4 (18%)
31	MA6	a	1515	31	19,26,27	1.01	2 (10%)	18,38,41	3.71	2 (11%)
31	7MG	a	524	31	22,26,27	1.22	1 (4%)	29,39,42	0.81	2 (6%)
5	2MA	A	2499	5	17,25,26	2.54	5 (29%)	17,37,40	1.35	2 (11%)
31	MA6	a	1516	31	19,26,27	1.04	2 (10%)	18,38,41	3.69	2 (11%)
52	H2U	v	20	52	18,21,22	0.56	0	21,30,33	1.12	1 (4%)
5	7MG	A	2065	5	22,26,27	1.19	1 (4%)	29,39,42	0.83	2 (6%)
31	2MG	a	1204	31	18,26,27	1.14	2 (11%)	16,38,41	0.94	1 (6%)
31	5MC	a	964	31	18,22,23	0.57	0	26,32,35	0.59	0
5	OMU	A	2548	5	19,22,23	3.06	8 (42%)	26,31,34	1.68	5 (19%)
5	PSU	A	2453	5	18,21,22	1.09	1 (5%)	22,30,33	1.82	6 (27%)
52	4SU	v	8	52	18,21,22	4.36	8 (44%)	26,30,33	2.32	8 (30%)
5	2MG	A	2441	5	18,26,27	1.14	2 (11%)	16,38,41	0.90	1 (6%)
5	PSU	A	952	5	18,21,22	1.08	1 (5%)	22,30,33	1.78	5 (22%)
31	UR3	a	1495	31	19,22,23	2.89	6 (31%)	26,32,35	1.31	2 (7%)
5	5MU	A	1935	5	19,22,23	0.43	0	28,32,35	0.53	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
52	PSU	v	56	52	-	2/7/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PSU	A	2576	5	-	0/7/25/26	0/2/2/2
52	5MU	v	55	52	-	3/7/25/26	0/2/2/2
31	2MG	a	963	31	-	2/5/27/28	0/3/3/3
31	PSU	a	513	31	-	2/7/25/26	0/2/2/2
5	PSU	A	2500	5	-	0/7/25/26	0/2/2/2
31	4OC	a	1399	31	-	2/9/29/30	0/2/2/2
5	6MZ	A	2026	5	-	1/5/27/28	0/3/3/3
5	OMG	A	2247	52,5	-	1/5/27/28	0/3/3/3
5	PSU	A	2601	5	-	0/7/25/26	0/2/2/2
31	MA6	a	1515	31	-	3/7/29/30	0/3/3/3
31	7MG	a	524	31	-	2/7/37/38	0/3/3/3
5	2MA	A	2499	5	-	3/3/25/26	0/3/3/3
31	MA6	a	1516	31	-	5/7/29/30	0/3/3/3
52	H2U	v	20	52	-	4/7/38/39	0/2/2/2
5	7MG	A	2065	5	-	1/7/37/38	0/3/3/3
31	2MG	a	1204	31	-	2/5/27/28	0/3/3/3
31	5MC	a	964	31	-	2/7/25/26	0/2/2/2
5	OMU	A	2548	5	-	2/9/27/28	0/2/2/2
5	PSU	A	2453	5	-	0/7/25/26	0/2/2/2
52	4SU	v	8	52	-	5/7/25/26	0/2/2/2
5	2MG	A	2441	5	-	0/5/27/28	0/3/3/3
5	PSU	A	952	5	-	0/7/25/26	0/2/2/2
31	UR3	a	1495	31	-	0/7/25/26	0/2/2/2
5	5MU	A	1935	5	-	0/7/25/26	0/2/2/2

The worst 5 of 59 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
52	v	8	4SU	C2-N1	10.04	1.54	1.38
52	v	8	4SU	C4-N3	8.38	1.46	1.37
5	A	2499	2MA	C2-N3	7.39	1.46	1.31
5	A	2548	OMU	C2-N1	7.19	1.50	1.38
31	a	1495	UR3	C2-N1	7.04	1.48	1.38

The worst 5 of 69 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	a	1515	MA6	N1-C6-N6	-14.49	101.81	117.06
31	a	1516	MA6	N1-C6-N6	-14.41	101.89	117.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2026	6MZ	C1'-N9-C4	-11.40	106.61	126.64
52	v	8	4SU	C4-N3-C2	-7.12	120.42	127.34
31	a	1515	MA6	N3-C2-N1	-5.54	120.02	128.68

There are no chirality outliers.

5 of 42 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
31	a	513	PSU	C3'-C4'-C5'-O5'
31	a	513	PSU	O4'-C4'-C5'-O5'
31	a	524	7MG	C3'-C4'-C5'-O5'
31	a	963	2MG	O4'-C4'-C5'-O5'
31	a	1399	4OC	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
55	80P	a	1603	56	44,45,45	3.53	21 (47%)	52,72,72	2.46	17 (32%)
55	80P	a	1601	56	44,45,45	3.40	20 (45%)	52,72,72	2.45	15 (28%)
55	80P	a	1602	56	44,45,45	3.82	22 (50%)	52,72,72	2.47	18 (34%)

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
55	80P	a	1603	56	-	13/22/87/87	0/5/5/5
55	80P	a	1601	56	-	12/22/87/87	0/5/5/5
55	80P	a	1602	56	-	12/22/87/87	0/5/5/5

The worst 5 of 63 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
55	a	1602	80P	C08-C09	13.07	1.63	1.52
55	a	1603	80P	C08-C09	11.93	1.62	1.52
55	a	1601	80P	C08-C09	11.75	1.62	1.52
55	a	1602	80P	C08-C21	9.94	1.62	1.53
55	a	1601	80P	C08-C21	8.34	1.60	1.53

The worst 5 of 50 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
55	a	1603	80P	C11-C10-C09	8.82	125.77	118.80
55	a	1601	80P	C11-C10-C09	8.68	125.66	118.80
55	a	1602	80P	C11-C10-C09	8.44	125.47	118.80
55	a	1601	80P	C08-C21-C04	5.17	118.71	111.64
55	a	1602	80P	C12-C11-C10	5.15	127.30	118.53

There are no chirality outliers.

5 of 37 torsion outliers are listed below:

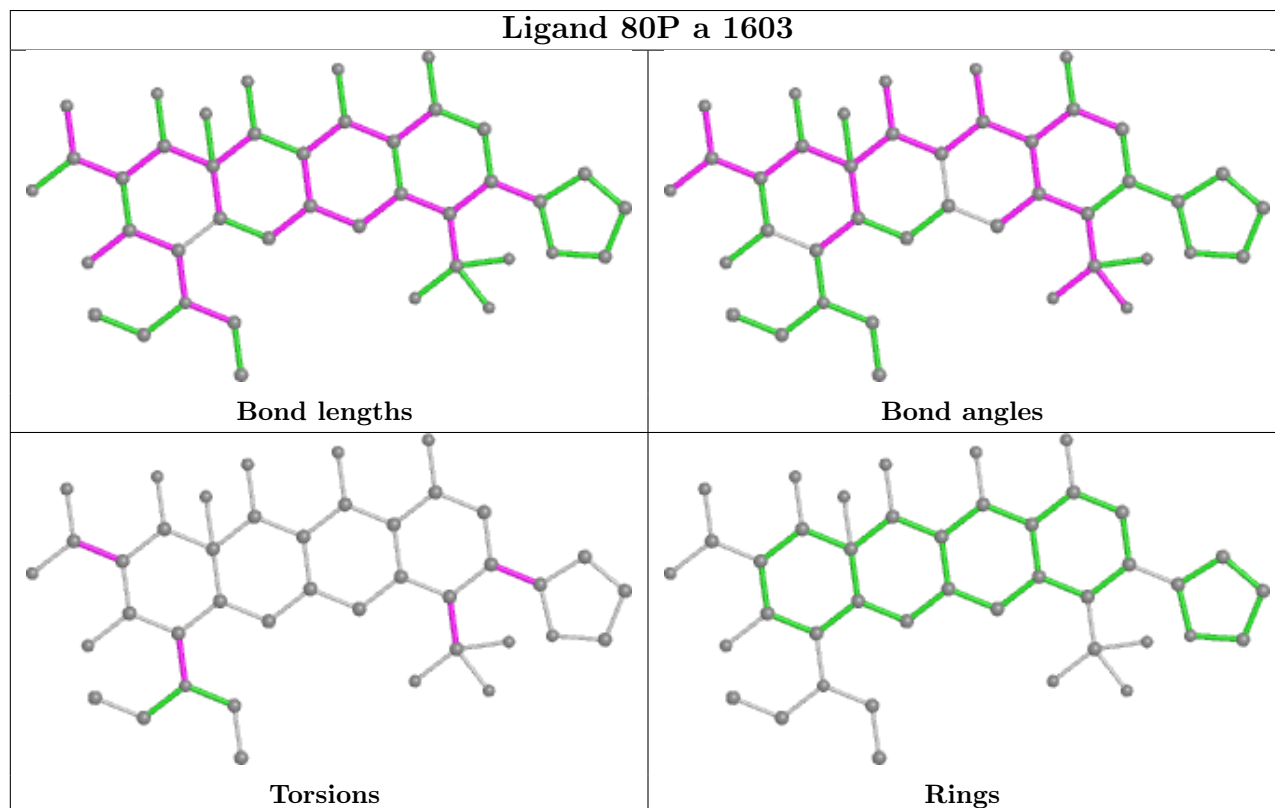
Mol	Chain	Res	Type	Atoms
55	a	1601	80P	C05-C04-N03-C01
55	a	1601	80P	C21-C04-N03-C40
55	a	1601	80P	C14-C15-C26-N27
55	a	1601	80P	C16-C15-C26-C30
55	a	1601	80P	C16-C15-C26-N27

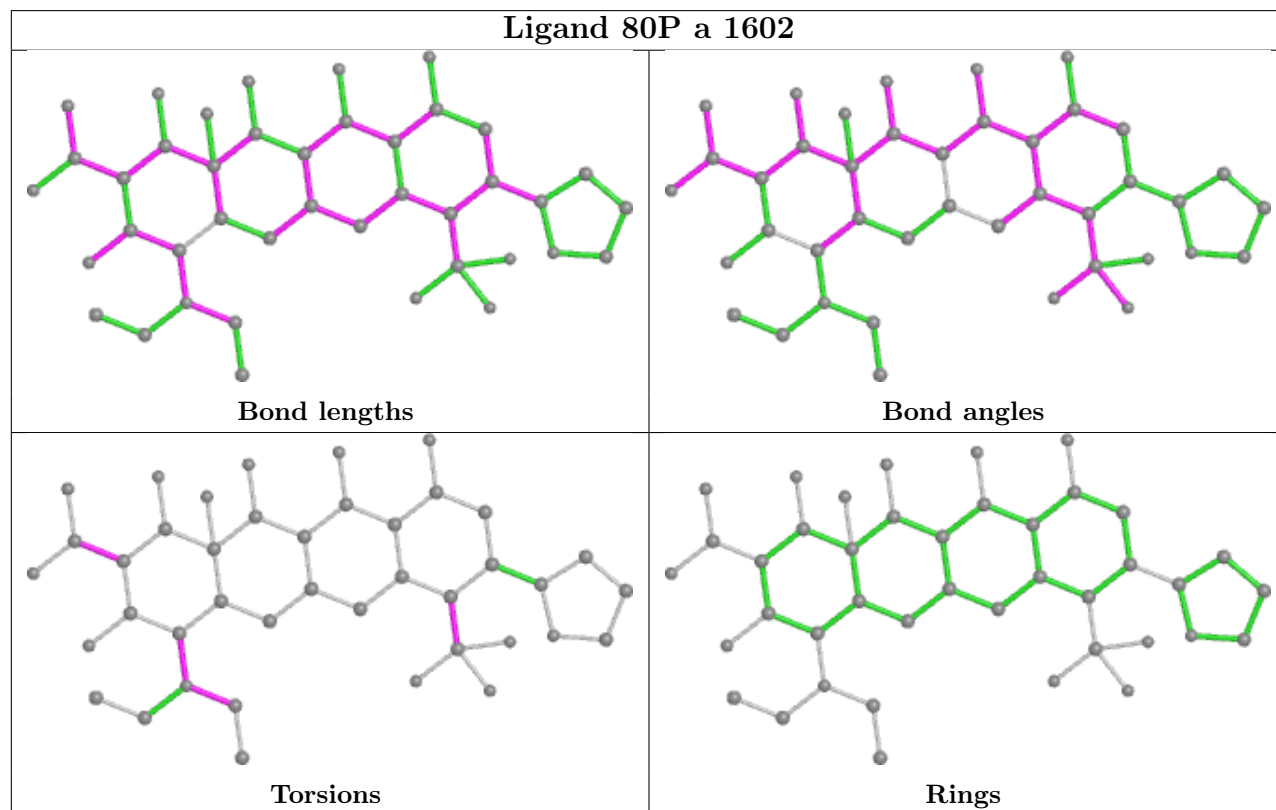
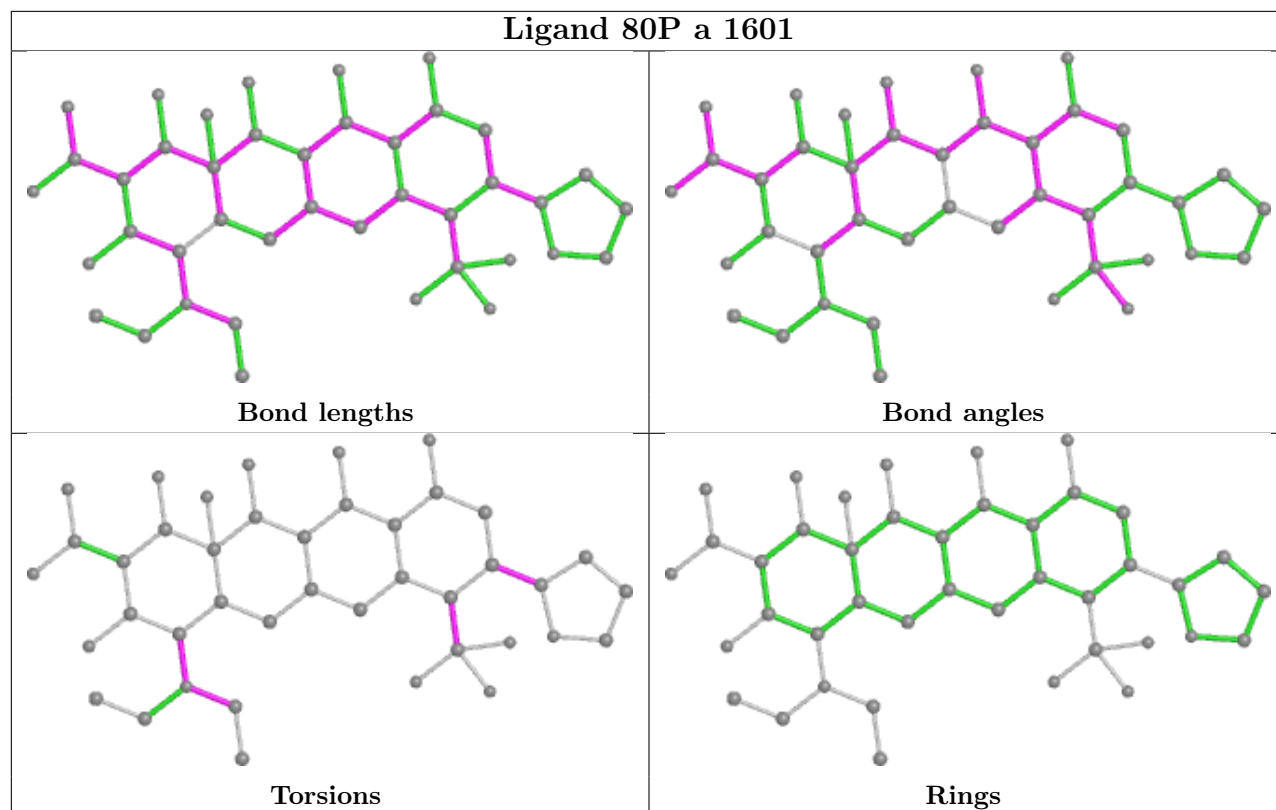
There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

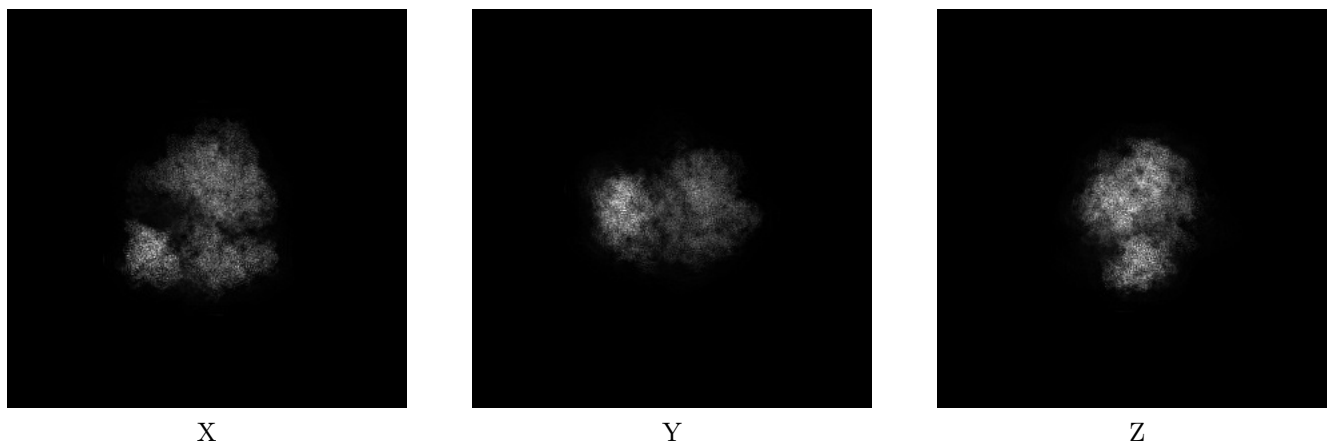
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24738. These allow visual inspection of the internal detail of the map and identification of artifacts.

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

6.1 Orthogonal projections [i](#)

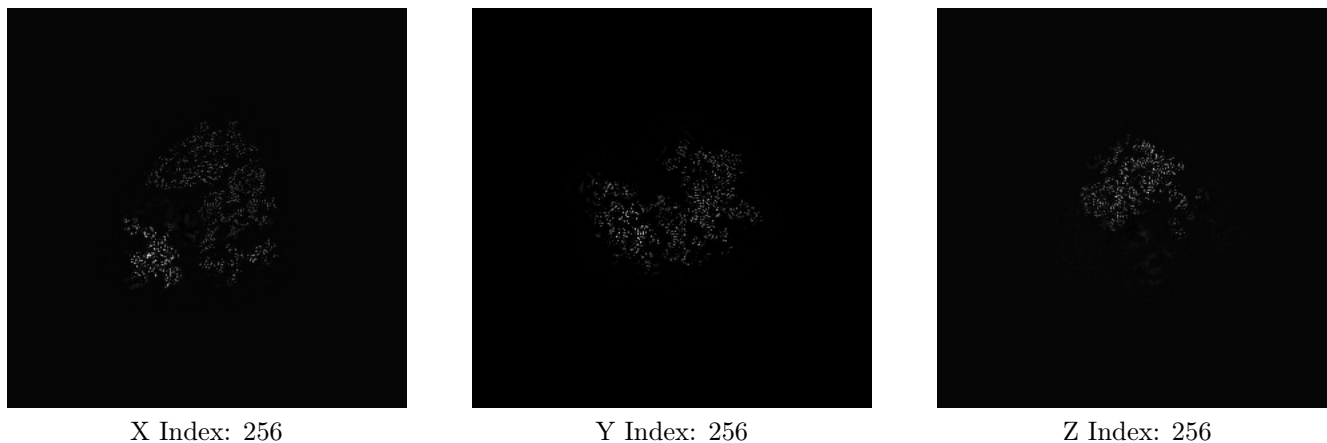
6.1.1 Primary map



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

6.2 Central slices [i](#)

6.2.1 Primary map



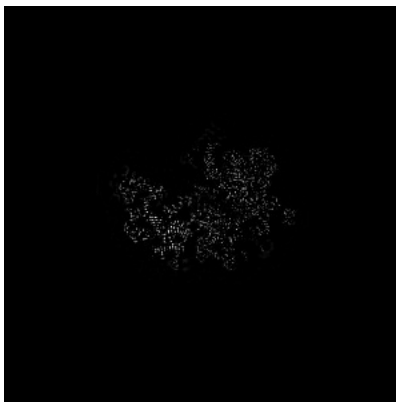
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

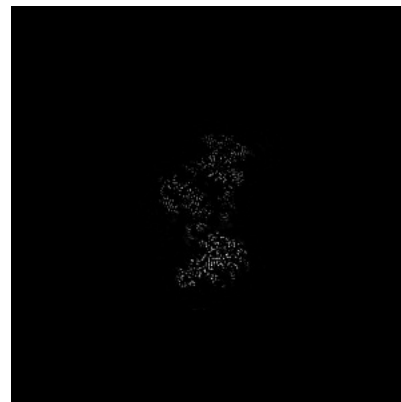
6.3.1 Primary map



X Index: 263



Y Index: 265



Z Index: 194

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

6.4 Orthogonal surface views [i](#)

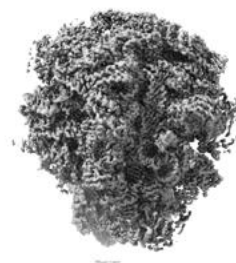
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

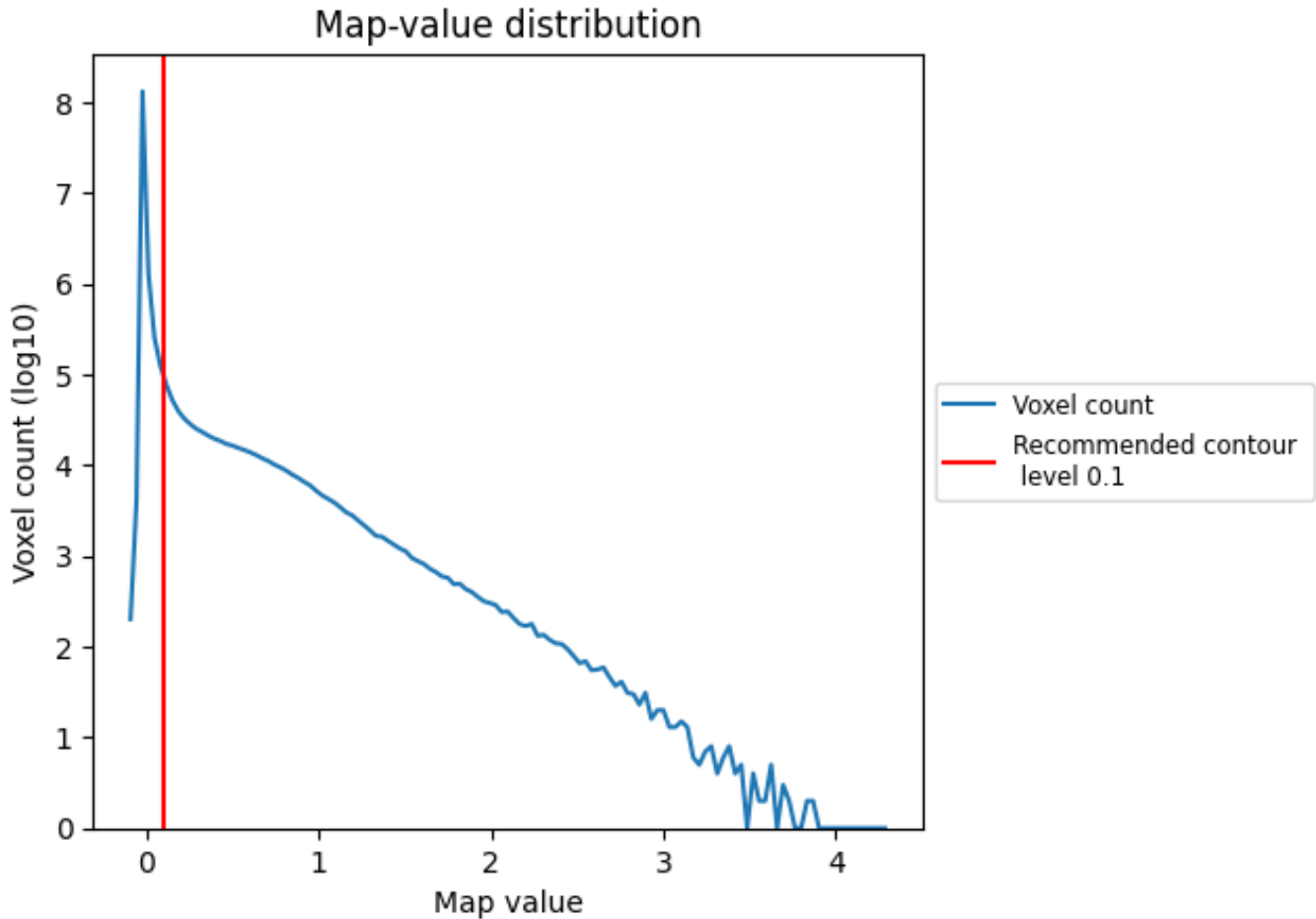
6.5 Mask visualisation

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

7 Map analysis [i](#)

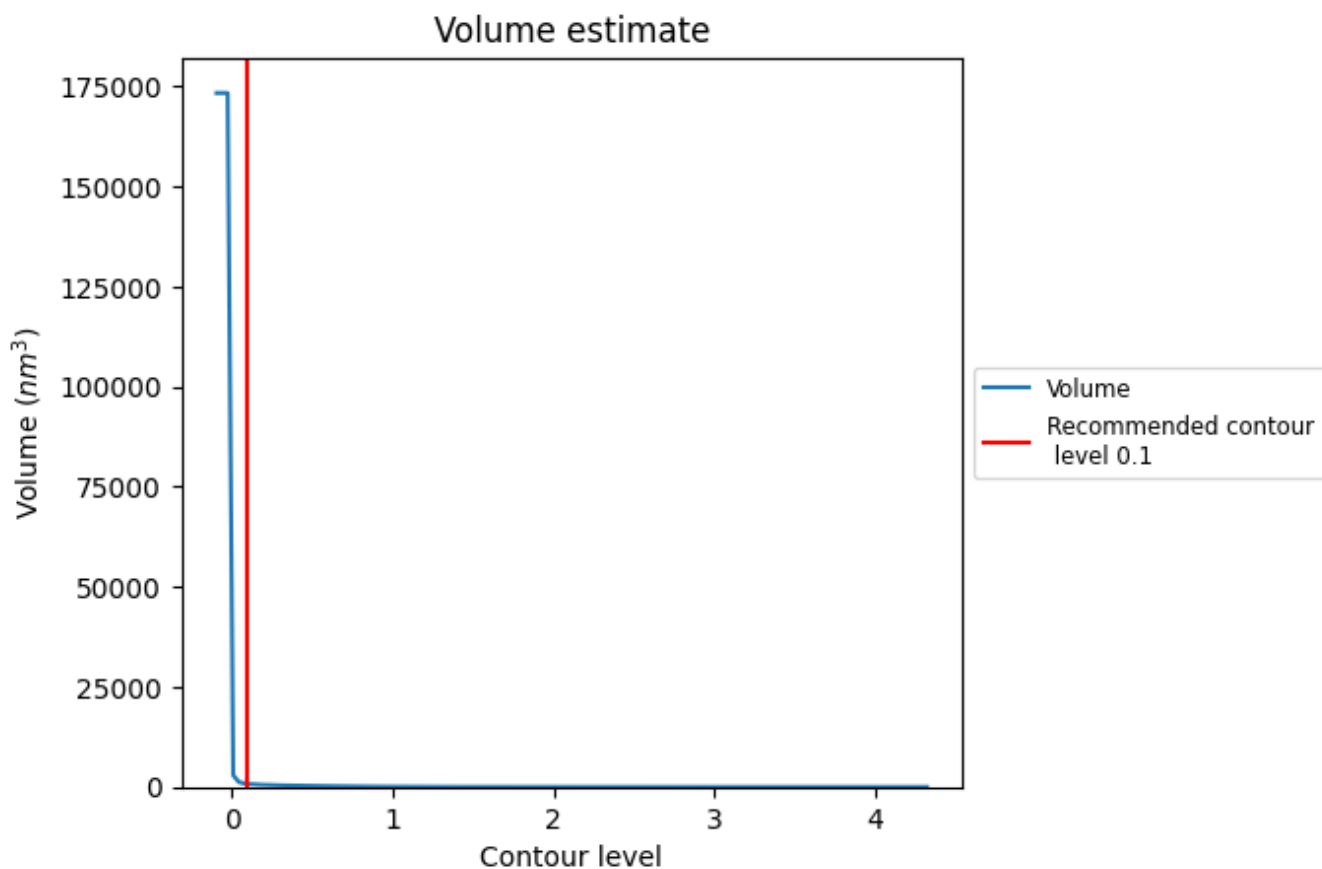
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

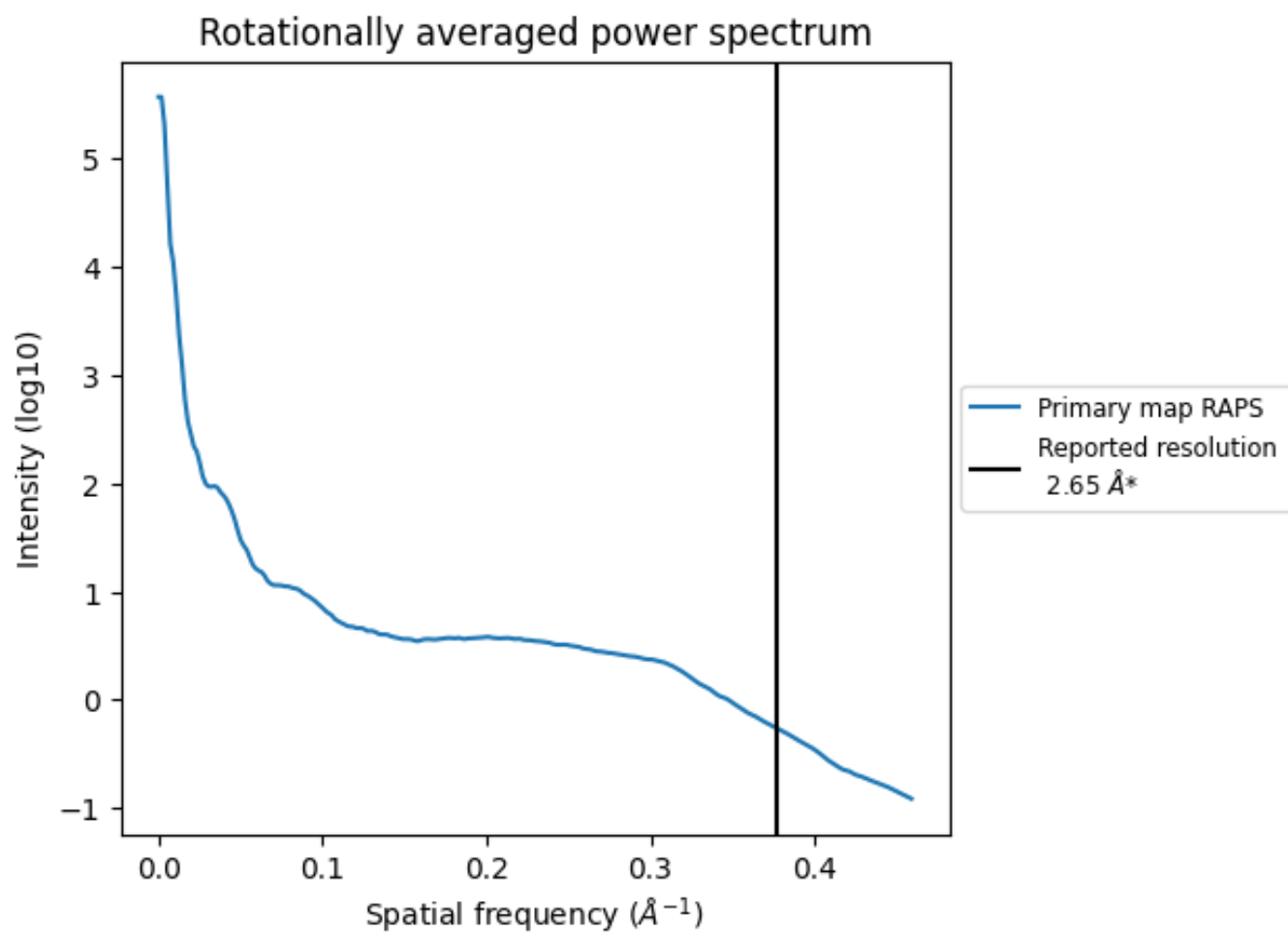
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 789 nm³; this corresponds to an approximate mass of 713 kDa.

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

7.3 Rotationally averaged power spectrum [i](#)



*Reported resolution corresponds to spatial frequency of 0.377 Å⁻¹

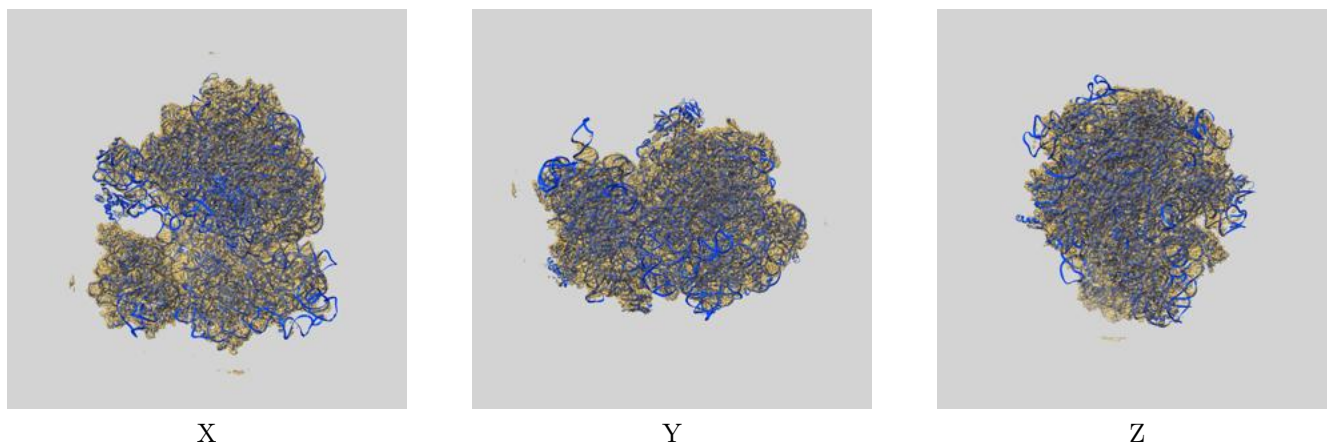
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

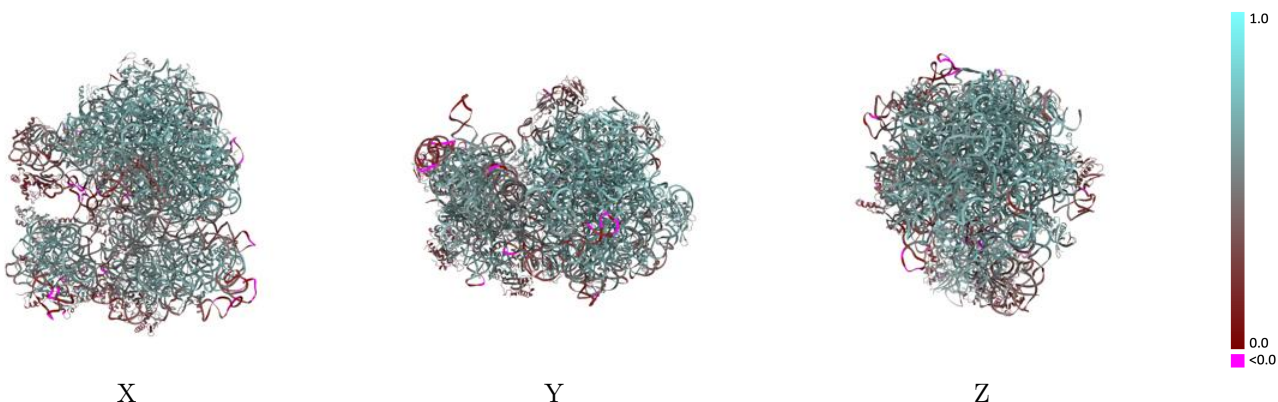
This section contains information regarding the fit between EMDB map EMD-24738 and PDB model 7RYF. Per-residue inclusion information can be found in section 3 on page 14.

9.1 Map-model overlay [i](#)



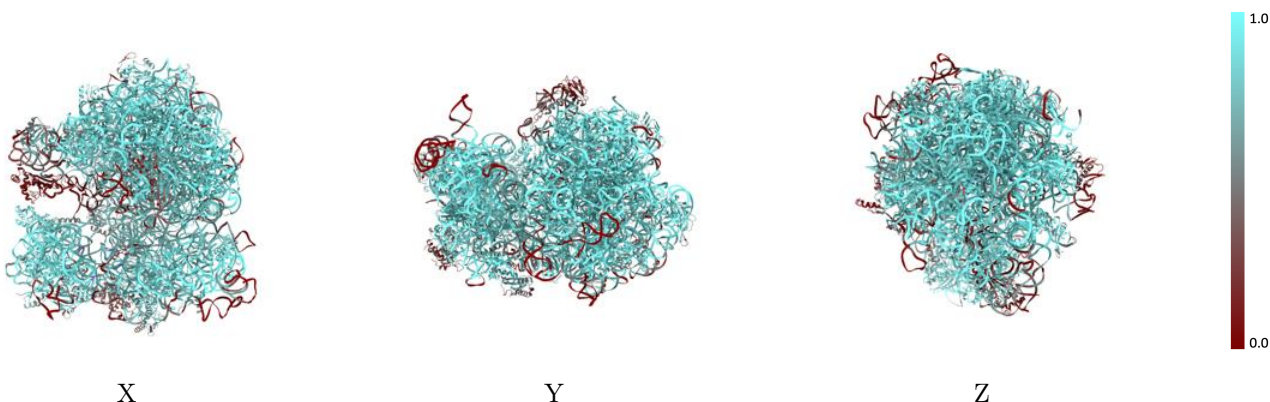
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



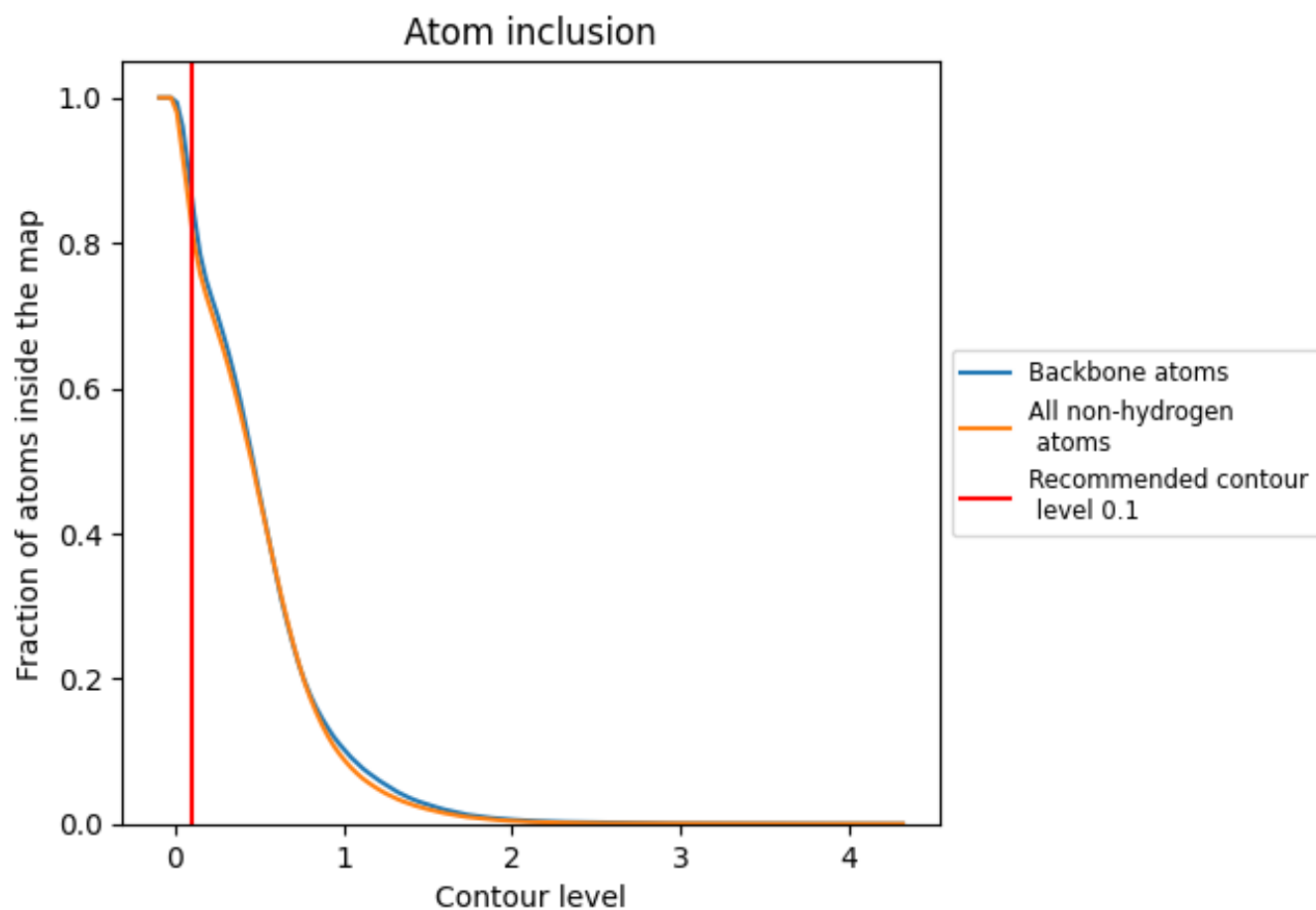
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.1).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 82% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8183	 0.5410
0	 0.7692	 0.5310
1	 0.9708	 0.6830
2	 0.8689	 0.5990
3	 0.6469	 0.4940
A	 0.8720	 0.5910
B	 0.5653	 0.4400
C	 0.9208	 0.6280
D	 0.9380	 0.6570
E	 0.8596	 0.6160
F	 0.0689	 0.2260
G	 0.2757	 0.3480
H	 0.3541	 0.3450
I	 0.9372	 0.6580
J	 0.9151	 0.6250
K	 0.9383	 0.6450
L	 0.8574	 0.5830
M	 0.9714	 0.6830
N	 0.5899	 0.4030
O	 0.8670	 0.5880
P	 0.9731	 0.6920
Q	 0.9048	 0.6270
R	 0.9258	 0.6570
S	 0.7696	 0.5630
T	 0.7103	 0.5410
U	 0.7995	 0.5550
V	 0.9136	 0.6500
W	 0.8858	 0.6140
X	 0.5903	 0.4520
Y	 0.8985	 0.6240
Z	 0.8756	 0.6180
a	 0.8404	 0.5080
b	 0.2733	 0.2740
c	 0.8860	 0.5030
d	 0.5552	 0.3520



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Chain	Atom inclusion	Q-score
e	 0.8976	 0.5460
f	 0.5623	 0.3430
g	 0.6951	 0.3510
h	 0.9441	 0.5970
i	 0.9517	 0.5770
j	 0.8729	 0.5130
k	 0.5113	 0.3240
l	 0.6487	 0.4530
m	 0.9171	 0.5150
n	 0.9294	 0.5680
o	 0.9063	 0.5580
p	 0.9084	 0.5970
q	 0.8433	 0.5070
r	 0.7869	 0.4540
s	 0.9556	 0.5670
t	 0.9174	 0.5620
u	 0.1722	 0.2610
v	 0.2934	 0.1800
w	 0.7692	 0.3290