



wwPDB EM Validation Summary Report ⓘ

Mar 10, 2026 – 04:24 PM UTC

PDB ID : 7QI5 / pdb_00007qi5
EMDB ID : EMD-13981
Title : Human mitochondrial ribosome in complex with mRNA, A/A-, P/P- and E/E-tRNAs at 2.63 Å resolution
Authors : Singh, V.; Itoh, Y.; Amunts, A.
Deposited on : 2021-12-14
Resolution : 2.63 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

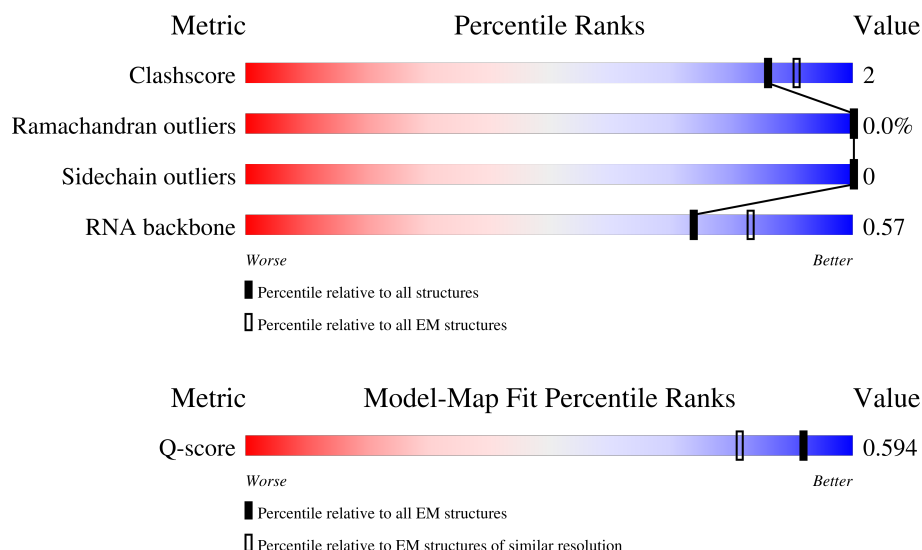
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.63 Å.

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



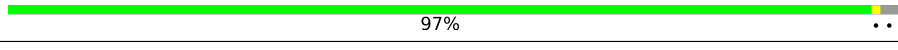
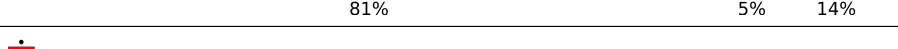
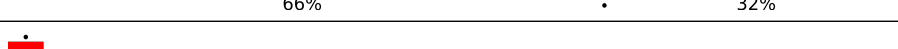
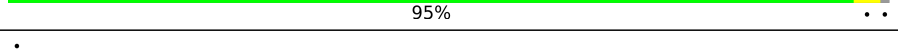
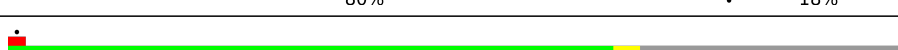
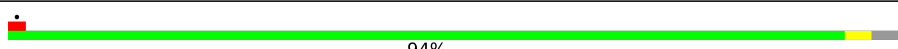
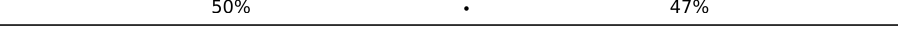
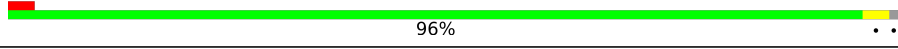
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	217187	24013	-
Ramachandran outliers	211220	23611	-
Sidechain outliers	210688	23127	-
RNA backbone	8275	3510	-
Q-score	-	25397	8888 (2.13 - 3.13)

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	AA	954	
2	AB	296	
3	AC	167	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	AD	430	
5	AE	125	
6	AF	242	
7	AG	396	
8	AH	201	
9	AI	194	
10	AJ	138	
11	AK	128	
12	AL	257	
13	AM	137	
14	AN	130	
15	AO	258	
16	AP	142	
17	AQ	87	
18	AR	360	
19	AS	190	
20	AT	173	
21	AU	205	
22	AV	414	
23	AW	187	
24	AX	398	
25	AY	395	
26	AZ	106	
27	A0	217	
28	A1	323	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
29	A2	118	
30	A3	199	
31	A4	689	
32	Aw	68	
33	Ax	70	
34	Ay	70	
35	Az	34	
36	A	1561	
37	B	72	
38	D	305	
39	E	348	
40	F	311	
41	I	261	
42	J	192	
43	K	178	
44	L	145	
45	M	296	
46	N	251	
47	O	175	
48	P	180	
49	Q	292	
50	R	149	
51	S	205	
52	T	206	
53	U	153	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
54	V	216	
55	W	148	
56	X	256	
57	Y	250	
58	Z	161	
59	0	188	
60	1	65	
61	2	92	
62	3	188	
63	4	103	
64	5	423	
65	6	380	
66	7	338	
67	8	206	
68	9	137	
69	a	142	
70	b	215	
71	c	332	
72	d	306	
73	e	279	
74	f	212	
75	g	166	
76	h	158	
77	i	128	
78	j	123	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
79	k	112	
80	l	138	
81	m	128	
82	o	102	
83	p	206	
84	q	222	
85	r	196	
86	s	439	
87	t	198	
87	u	198	
87	v	198	
87	w	198	
87	x	198	
87	y	198	
88	H	267	
89	z	325	

2 Entry composition

There are 101 unique types of molecules in this entry. The entry contains 343544 atoms, of which 153978 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 12S mitochondrial rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	AA	954	Total	C	H	N	O	P	0	0
			30564	9088	10304	3647	6571	954		

- Molecule 2 is a protein called 28S ribosomal protein S2, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	AB	224	Total	C	H	N	O	S	0	0
			3627	1158	1809	328	322	10		

- Molecule 3 is a protein called 28S ribosomal protein S24, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	AC	132	Total	C	H	N	O	S	0	0
			2172	699	1089	195	185	4		

- Molecule 4 is a protein called 28S ribosomal protein S5, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	AD	343	Total	C	H	N	O	S	0	0
			5536	1713	2805	518	487	13		

- Molecule 5 is a protein called 28S ribosomal protein S6, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	AE	122	Total	C	H	N	O	S	0	0
			1972	614	1000	177	177	4		

- Molecule 6 is a protein called 28S ribosomal protein S7, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	AF	208	Total	C	H	N	O	S	0	0
			3496	1104	1771	312	298	11		

- Molecule 7 is a protein called 28S ribosomal protein S9, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	AG	327	Total	C	H	N	O	S	0	0
			5377	1710	2689	477	487	14		

- Molecule 8 is a protein called 28S ribosomal protein S10, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	AH	140	Total	C	H	N	O	S	0	0
			2339	745	1187	194	210	3		

- Molecule 9 is a protein called 28S ribosomal protein S11, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	AI	137	Total	C	H	N	O	S	0	0
			2081	642	1061	192	182	4		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AI	184	5F0	ASN	conflict	UNP P82912

- Molecule 10 is a protein called 28S ribosomal protein S12, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	AJ	108	Total	C	H	N	O	S	0	0
			1728	521	889	169	143	6		

- Molecule 11 is a protein called 28S ribosomal protein S14, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
11	AK	101	Total	C	H	N	O	S	0	0
			1748	537	886	179	141	5		

- Molecule 12 is a protein called 28S ribosomal protein S15, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
12	AL	174	Total	C	H	N	O	S	0	0
			2994	925	1541	270	251	7		

- Molecule 13 is a protein called 28S ribosomal protein S16, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
13	AM	119	Total	C	H	N	O	S	0	0
			1908	594	966	185	157	6		

- Molecule 14 is a protein called 28S ribosomal protein S17, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
14	AN	110	Total	C	H	N	O	S	0	0
			1797	562	929	156	147	3		

- Molecule 15 is a protein called 28S ribosomal protein S18b, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
15	AO	193	Total	C	H	N	O	S	0	0
			3153	1014	1561	294	277	7		

- Molecule 16 is a protein called 28S ribosomal protein S18c, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
16	AP	97	Total	C	H	N	O	S	0	0
			1588	501	807	134	138	8		

- Molecule 17 is a protein called MRPS21 isoform 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
17	AQ	86	Total	C	H	N	O	S	0	0
			1502	460	758	150	126	8		

- Molecule 18 is a protein called 28S ribosomal protein S22, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
18	AR	295	Total	C	H	N	O	S	0	0
			4839	1533	2430	413	455	8		

- Molecule 19 is a protein called 28S ribosomal protein S23, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
19	AS	135	Total	C	H	N	O	S	0	0
			2227	716	1116	198	196	1		

- Molecule 20 is a protein called 28S ribosomal protein S25, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
20	AT	168	Total	C	H	N	O	S	0	0
			2764	877	1393	239	244	11		

- Molecule 21 is a protein called 28S ribosomal protein S26, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
21	AU	176	Total	C	H	N	O	S	0	0
			2989	916	1501	301	267	4		

- Molecule 22 is a protein called 28S ribosomal protein S27, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
22	AV	362	Total	C	H	N	O	S	0	0
			5933	1904	2964	495	558	12		

- Molecule 23 is a protein called 28S ribosomal protein S28, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
23	AW	100	Total	C	H	N	O	S	0	0
			1593	498	804	141	146	4		

- Molecule 24 is a protein called 28S ribosomal protein S29, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
24	AX	352	Total	C	H	N	O	S	0	0
			5694	1822	2845	499	517	11		

- Molecule 25 is a protein called 28S ribosomal protein S31, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
25	AY	149	Total	C	H	N	O	S	0	0
			2444	801	1198	207	234	4		

- Molecule 26 is a protein called 28S ribosomal protein S33, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
26	AZ	100	Total	C	H	N	O	S	0	0
			1698	534	859	153	148	4		

- Molecule 27 is a protein called 28S ribosomal protein S34, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
27	A0	215	Total	C	H	N	O	S	0	0
			3584	1130	1797	339	313	5		

- Molecule 28 is a protein called 28S ribosomal protein S35, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
28	A1	279	Total	C	H	N	O	S	0	0
			4561	1435	2296	387	432	11		

- Molecule 29 is a protein called Coiled-coil-helix-coiled-coil-helix domain-containing protein 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
29	A2	117	Total	C	H	N	O	S	0	0
			1906	579	971	182	166	8		

- Molecule 30 is a protein called Aurora kinase A-interacting protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
30	A3	70	Total	C	H	N	O	S	0	0
			1326	401	701	134	89	1		

- Molecule 31 is a protein called Pentatricopeptide repeat domain-containing protein 3, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
31	A4	588	Total	C	H	N	O	S	0	0
			9538	3053	4770	808	879	28		

- Molecule 32 is a RNA chain called A/A-tRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
32	Aw	68	Total	C	H	N	O	P	0	0
			2159	646	725	248	472	68		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Aw	26	U	G	conflict	GB 1896813690
Aw	34	A	U	conflict	GB 1896813690
Aw	35	A	C	conflict	GB 1896813690
Aw	36	C	G	conflict	GB 1896813690
Aw	73	U	A	conflict	GB 1896813690

- Molecule 33 is a RNA chain called P/P-tRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
33	Ax	70	Total	C	H	N	O	P	0	0
			2233	665	751	260	487	70		

- Molecule 34 is a RNA chain called E/E-tRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
34	Ay	70	Total	C	H	N	O	P	0	0
			2235	665	752	261	487	70		

- Molecule 35 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
35	Az	34	Total	C	H	N	O	P	0	0
			1079	324	360	123	238	34		

- Molecule 36 is a RNA chain called 16S mitochondrial rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
36	A	1558	Total	C	H	N	O	P	0	0
			49871	14843	16801	5963	10706	1558		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3107	U	C	conflict	GB 208964619

- Molecule 37 is a RNA chain called mitochondrial tRNAVal.

Mol	Chain	Residues	Atoms						AltConf	Trace
37	B	72	Total	C	H	N	O	P	0	0
			2303	685	779	269	498	72		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	74	C	G	conflict	GB NC_012920.1
B	76	A	U	conflict	GB NC_012920.1

- Molecule 38 is a protein called 39S ribosomal protein L2, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
38	D	238	Total	C	H	N	O	S	0	0
			3780	1157	1921	376	317	9		

- Molecule 39 is a protein called 39S ribosomal protein L3, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
39	E	305	Total	C	H	N	O	S	0	0
			4822	1545	2416	418	432	11		

- Molecule 40 is a protein called 39S ribosomal protein L4, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
40	F	252	Total	C	H	N	O	S	0	0
			4097	1305	2066	370	350	6		

- Molecule 41 is a protein called 39S ribosomal protein L10, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
41	I	212	Total	C	H	N	O	S	0	0
			3484	1088	1789	304	292	11		

- Molecule 42 is a protein called 39S ribosomal protein L11, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
42	J	175	Total	C	H	N	O	S	0	0
			2739	847	1409	237	244	2		

- Molecule 43 is a protein called 39S ribosomal protein L13, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
43	K	177	Total	C	H	N	O	S	0	0
			2907	936	1452	259	253	7		

- Molecule 44 is a protein called 39S ribosomal protein L14, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
44	L	115	Total	C	H	N	O	S	0	0
			1832	559	942	171	155	5		

- Molecule 45 is a protein called 39S ribosomal protein L15, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
45	M	289	Total	C	H	N	O	S	0	0
			4699	1476	2385	427	405	6		

- Molecule 46 is a protein called 39S ribosomal protein L16, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
46	N	222	Total	C	H	N	O	S	0	0
			3604	1143	1818	326	307	10		

- Molecule 47 is a protein called 39S ribosomal protein L17, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
47	O	154	Total	C	H	N	O	S	0	0
			2554	792	1295	241	219	7		

- Molecule 48 is a protein called 39S ribosomal protein L18, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
48	P	144	Total	C	H	N	O	S	0	0
			2339	733	1166	224	211	5		

- Molecule 49 is a protein called 39S ribosomal protein L19, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
49	Q	239	Total	C	H	N	O	S	0	0
			4021	1277	2031	353	351	9		

- Molecule 50 is a protein called 39S ribosomal protein L20, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
50	R	140	Total	C	H	N	O	S	0	0
			2369	732	1215	231	187	4		

- Molecule 51 is a protein called 39S ribosomal protein L21, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
51	S	161	Total	C	H	N	O	S	0	0
			2659	835	1366	227	227	4		

- Molecule 52 is a protein called 39S ribosomal protein L22, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
52	T	166	Total	C	H	N	O	S	0	0
			2781	875	1412	254	233	7		

- Molecule 53 is a protein called 39S ribosomal protein L23, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
53	U	152	Total	C	H	N	O	S	0	0
			2483	788	1232	234	226	3		

- Molecule 54 is a protein called 39S ribosomal protein L24, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
54	V	205	Total	C	H	N	O	S	0	0
			3365	1068	1689	298	302	8		

- Molecule 55 is a protein called 39S ribosomal protein L27, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
55	W	116	Total	C	H	N	O	S	0	0
			1840	577	936	171	153	3		

- Molecule 56 is a protein called 39S ribosomal protein L28, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
56	X	244	Total	C	H	N	O	S	0	0
			4105	1322	2061	352	365	5		

- Molecule 57 is a protein called 39S ribosomal protein L47, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
57	Y	181	Total	C	H	N	O	S	0	0
			3154	995	1598	298	259	4		

- Molecule 58 is a protein called 39S ribosomal protein L30, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
58	Z	122	Total	C	H	N	O	S	0	0
			2041	636	1045	186	171	3		

- Molecule 59 is a protein called 39S ribosomal protein L32, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
59	0	110	Total	C	H	N	O	S	0	0
			1815	554	917	176	162	6		

- Molecule 60 is a protein called 39S ribosomal protein L33, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
60	1	56	Total	C	H	N	O	S	0	0
			977	296	513	89	77	2		

- Molecule 61 is a protein called 39S ribosomal protein L34, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
61	2	46	Total	C	H	N	O	S	0	0
			784	233	407	83	60	1		

- Molecule 62 is a protein called 39S ribosomal protein L35, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
62	3	95	Total	C	H	N	O	S	0	0
			1716	539	884	162	128	3		

- Molecule 63 is a protein called 39S ribosomal protein L36, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
63	4	38	Total	C	H	N	O	S	0	0
			704	217	362	72	49	4		

- Molecule 64 is a protein called 39S ribosomal protein L37, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
64	5	394	Total	C	H	N	O	S	0	0
			6419	2073	3209	560	566	11		

- Molecule 65 is a protein called 39S ribosomal protein L38, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
65	6	354	Total	C	H	N	O	S	0	0
			5792	1881	2844	525	533	9		

- Molecule 66 is a protein called 39S ribosomal protein L39, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
66	7	294	Total	C	H	N	O	S	0	0
			4789	1529	2399	405	438	18		

- Molecule 67 is a protein called 39S ribosomal protein L40, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
67	8	157	Total	C	H	N	O	S	0	0
			2696	844	1369	235	246	2		

- Molecule 68 is a protein called 39S ribosomal protein L41, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
68	9	124	Total	C	H	N	O	S	0	0
			1985	644	988	170	181	2		

- Molecule 69 is a protein called 39S ribosomal protein L42, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
69	a	100	Total	C	H	N	O	S	0	0
			1652	529	812	152	154	5		

- Molecule 70 is a protein called Large ribosomal subunit protein mL43.

Mol	Chain	Residues	Atoms						AltConf	Trace
70	b	151	Total	C	H	N	O	S	0	0
			2392	744	1196	231	218	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
b	2	ACE	-	acetylation	UNP Q8N983

- Molecule 71 is a protein called 39S ribosomal protein L44, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
71	c	286	Total	C	H	N	O	S	0	0
			4621	1470	2322	397	423	9		

- Molecule 72 is a protein called 39S ribosomal protein L45, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
72	d	241	Total	C	H	N	O	S	0	0
			3964	1273	1979	340	359	13		

- Molecule 73 is a protein called 39S ribosomal protein L46, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
73	e	238	Total	C	H	N	O	S	0	0
			3848	1222	1917	339	364	6		

- Molecule 74 is a protein called 39S ribosomal protein L48, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
74	f	157	Total	C	H	N	O	S	0	0
			2523	799	1271	207	242	4		

- Molecule 75 is a protein called 39S ribosomal protein L49, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
75	g	134	Total	C	H	N	O	S	0	0
			2210	719	1097	193	199	2		

- Molecule 76 is a protein called 39S ribosomal protein L50, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
76	h	110	Total	C	H	N	O	S	0	0
			1777	568	882	156	168	3		

- Molecule 77 is a protein called 39S ribosomal protein L51, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
77	i	97	Total	C	H	N	O	S	0	0
			1687	532	859	165	127	4		

- Molecule 78 is a protein called 39S ribosomal protein L52, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
78	j	94	Total	C	H	N	O	S	0	0
			1492	463	747	144	136	2		

- Molecule 79 is a protein called 39S ribosomal protein L53, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
79	k	101	Total	C	H	N	O	S	0	0
			1559	479	785	148	142	5		

- Molecule 80 is a protein called 39S ribosomal protein L54, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
80	l	82	Total	C	H	N	O	S	0	0
			1363	437	675	120	128	3		

- Molecule 81 is a protein called 39S ribosomal protein L55, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
81	m	92	Total	C	H	N	O	S	0	0
			1551	488	760	159	142	2		

- Molecule 82 is a protein called Ribosomal protein 63, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
82	o	94	Total	C	H	N	O	S	0	0
			1604	501	806	165	129	3		

- Molecule 83 is a protein called Peptidyl-tRNA hydrolase ICT1, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
83	p	147	Total	C	H	N	O	S	0	0
			2429	748	1224	228	225	4		

- Molecule 84 is a protein called Growth arrest and DNA damage-inducible proteins-interacting protein 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
84	q	165	Total	C	H	N	O	S	0	0
			2765	865	1376	270	249	5		

- Molecule 85 is a protein called 39S ribosomal protein S18a, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
85	r	162	Total	C	H	N	O	S	0	0
			2671	839	1349	252	223	8		

- Molecule 86 is a protein called 39S ribosomal protein S30, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
86	s	386	Total	C	H	N	O	S	0	0
			6298	2023	3143	559	559	14		

- Molecule 87 is a protein called 39S ribosomal protein L12, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
87	t	46	Total	C	H	N	O		0	0
			732	228	378	56	70			
87	u	32	Total	C	H	N	O		0	0
			541	168	284	40	49			
87	v	32	Total	C	H	N	O		0	0
			541	168	284	40	49			
87	w	31	Total	C	H	N	O		0	0
			520	159	275	39	47			
87	x	31	Total	C	H	N	O		0	0
			520	159	275	39	47			
87	y	31	Total	C	H	N	O		0	0
			520	159	275	39	47			

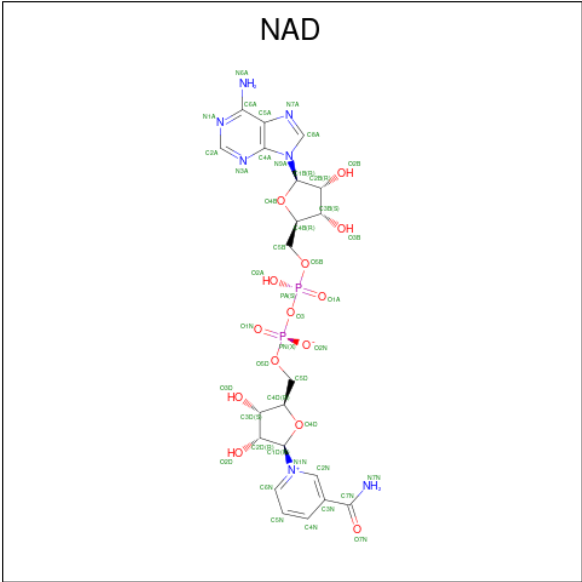
- Molecule 88 is a protein called 39S ribosomal protein L9, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
88	H	202	Total	C	H	N	O	S	0	0
			3397	1067	1736	304	286	4		

- Molecule 89 is a protein called 39S ribosomal protein L1, mitochondrial.

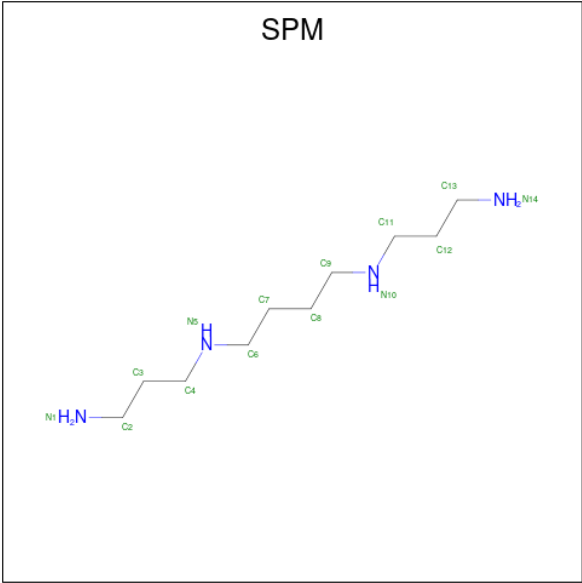
Mol	Chain	Residues	Atoms						AltConf	Trace
89	z	252	Total	C	H	N	O	S	0	0
			4104	1304	2077	336	381	6		

- Molecule 90 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



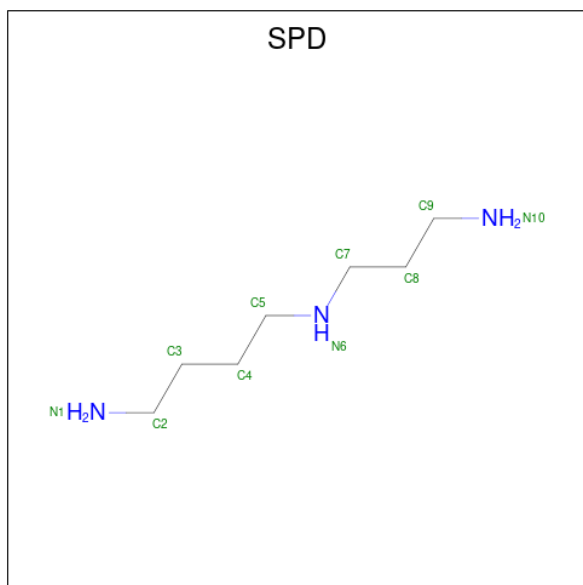
Mol	Chain	Residues	Atoms						AltConf
90	AA	1	Total	C	H	N	O	P	0
			70	21	26	7	14	2	

- Molecule 91 is SPERMINE (CCD ID: SPM) (formula: C₁₀H₂₆N₄).



Mol	Chain	Residues	Atoms				AltConf
91	AA	1	Total	C	H	N	0
			44	10	30	4	

- Molecule 92 is SPERMIDINE (CCD ID: SPD) (formula: C₇H₁₉N₃).



Mol	Chain	Residues	Atoms				AltConf
92	AA	1	Total	C	H	N	0
			32	7	22	3	
92	A	1	Total	C	H	N	0
			32	7	22	3	
92	A	1	Total	C	H	N	0
			32	7	22	3	
92	A	1	Total	C	H	N	0
			32	7	22	3	

- Molecule 93 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
93	AA	63	Total	Mg	0
			63	63	
93	AB	1	Total	Mg	0
			1	1	
93	AX	1	Total	Mg	0
			1	1	
93	A3	1	Total	Mg	0
			1	1	
93	A	138	Total	Mg	0
			138	138	
93	D	2	Total	Mg	0
			2	2	
93	E	1	Total	Mg	0
			1	1	
93	g	1	Total	Mg	0
			1	1	

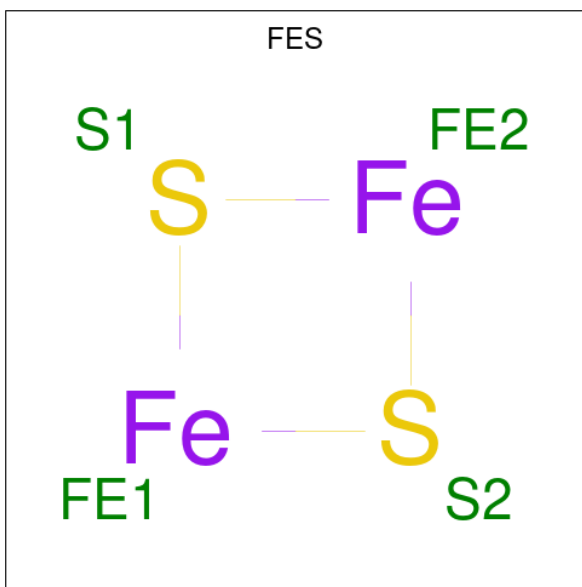
- Molecule 94 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
94	AA	18	Total 18	K 18	0
94	A	29	Total 29	K 29	0
94	D	1	Total 1	K 1	0
94	M	1	Total 1	K 1	0
94	W	1	Total 1	K 1	0
94	3	1	Total 1	K 1	0
94	6	1	Total 1	K 1	0
94	i	1	Total 1	K 1	0
94	o	1	Total 1	K 1	0

- Molecule 95 is ZINC ION (CCD ID: ZN) (formula: Zn).

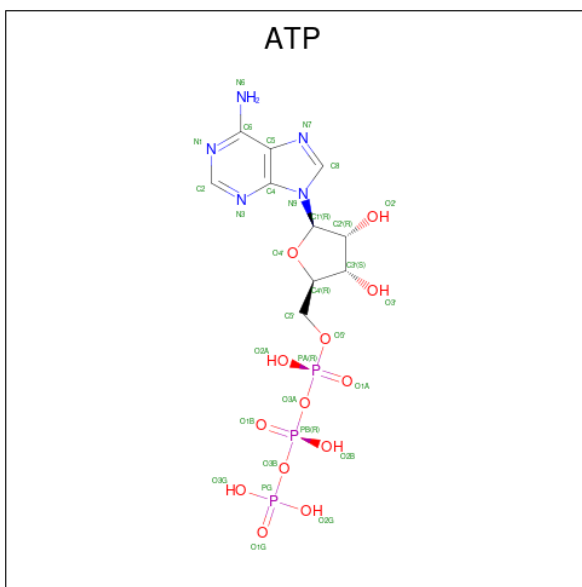
Mol	Chain	Residues	Atoms		AltConf
95	AO	1	Total 1	Zn 1	0
95	0	1	Total 1	Zn 1	0
95	4	1	Total 1	Zn 1	0

- Molecule 96 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			AltConf
96	AP	1	Total 4	Fe 2	S 2	0
96	AT	1	Total 4	Fe 2	S 2	0
96	r	1	Total 4	Fe 2	S 2	0

- Molecule 97 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



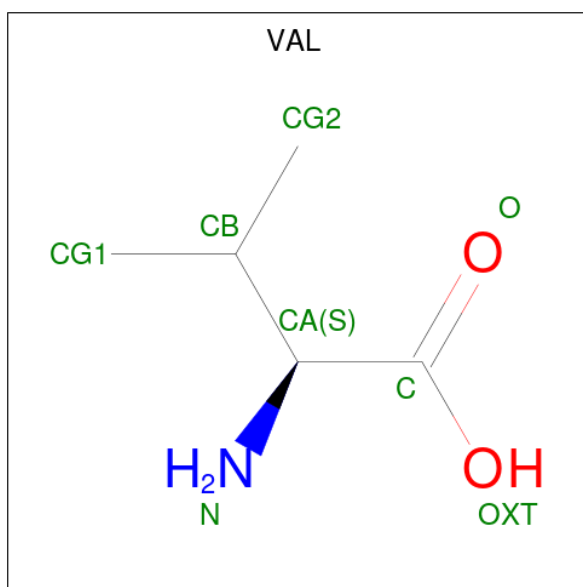
Mol	Chain	Residues	Atoms					AltConf	
97	AX	1	Total 43	C 10	H 12	N 5	O 13	P 3	0

-
- The image displays the chemical structure of GDP (Guanosine Diphosphate). It consists of a guanine base (a purine ring system with an amino group at C2 and a carbonyl group at C6) linked to a ribose sugar via a glycosidic bond at the N9 position. The ribose sugar is further linked to two phosphate groups (P1 and P2) via phosphodiester bonds. The structure is labeled with atom names and numbers, and the phosphate groups are shown with their respective negative charges (O⁻).

PUT

The diagram shows the chemical structure of Putrescine (PUT), a linear diamine. It consists of a four-carbon chain. The first carbon (C1) is bonded to an amino group (NH₂, N1). The second carbon (C2) is bonded to a hydrogen atom (H2). The third carbon (C3) is bonded to a hydrogen atom (H3). The fourth carbon (C4) is bonded to an amino group (NH₂, N2). The atoms are labeled with green text: C1, C2, C3, C4, N1, N2, H1, H2, H3, H4. The bonds are shown as black lines.





Mol	Chain	Residues	Atoms					AltConf
100	B	1	Total	C	H	N	O	0
			18	5	11	1	1	

- Molecule 101 is water.

Mol	Chain	Residues	Atoms		AltConf
101	AA	2039	Total	O	0
			2039	2039	
101	AB	78	Total	O	0
			78	78	
101	AC	55	Total	O	0
			55	55	
101	AD	82	Total	O	0
			82	82	
101	AE	27	Total	O	0
			27	27	
101	AF	39	Total	O	0
			39	39	
101	AG	64	Total	O	0
			64	64	
101	AH	57	Total	O	0
			57	57	
101	AI	39	Total	O	0
			39	39	
101	AJ	25	Total	O	0
			25	25	
101	AK	51	Total	O	0
			51	51	

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		AltConf
101	AL	42	Total 42	O 42	0
101	AM	22	Total 22	O 22	0
101	AN	39	Total 39	O 39	0
101	AO	37	Total 37	O 37	0
101	AP	31	Total 31	O 31	0
101	AQ	77	Total 77	O 77	0
101	AR	11	Total 11	O 11	0
101	AS	25	Total 25	O 25	0
101	AT	44	Total 44	O 44	0
101	AU	8	Total 8	O 8	0
101	AW	16	Total 16	O 16	0
101	AX	56	Total 56	O 56	0
101	AY	18	Total 18	O 18	0
101	AZ	30	Total 30	O 30	0
101	A0	1	Total 1	O 1	0
101	A1	35	Total 35	O 35	0
101	A2	32	Total 32	O 32	0
101	A3	50	Total 50	O 50	0
101	A4	9	Total 9	O 9	0
101	Aw	6	Total 6	O 6	0
101	Ax	7	Total 7	O 7	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		AltConf
101	Ay	3	Total 3	O 3	0
101	Az	13	Total 13	O 13	0
101	A	2955	Total 2955	O 2955	0
101	B	66	Total 66	O 66	0
101	D	68	Total 68	O 68	0
101	E	66	Total 66	O 66	0
101	F	79	Total 79	O 79	0
101	I	27	Total 27	O 27	0
101	J	1	Total 1	O 1	0
101	K	57	Total 57	O 57	0
101	L	31	Total 31	O 31	0
101	M	51	Total 51	O 51	0
101	N	58	Total 58	O 58	0
101	O	33	Total 33	O 33	0
101	P	81	Total 81	O 81	0
101	Q	29	Total 29	O 29	0
101	R	63	Total 63	O 63	0
101	S	49	Total 49	O 49	0
101	T	45	Total 45	O 45	0
101	U	24	Total 24	O 24	0
101	V	5	Total 5	O 5	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		AltConf
101	W	55	Total 55	O 55	0
101	X	11	Total 11	O 11	0
101	Y	25	Total 25	O 25	0
101	Z	32	Total 32	O 32	0
101	0	28	Total 28	O 28	0
101	1	3	Total 3	O 3	0
101	2	34	Total 34	O 34	0
101	3	44	Total 44	O 44	0
101	4	11	Total 11	O 11	0
101	5	14	Total 14	O 14	0
101	6	100	Total 100	O 100	0
101	7	15	Total 15	O 15	0
101	8	15	Total 15	O 15	0
101	9	17	Total 17	O 17	0
101	a	11	Total 11	O 11	0
101	b	35	Total 35	O 35	0
101	c	18	Total 18	O 18	0
101	d	9	Total 9	O 9	0
101	e	13	Total 13	O 13	0
101	f	19	Total 19	O 19	0
101	g	13	Total 13	O 13	0

Continued on next page...

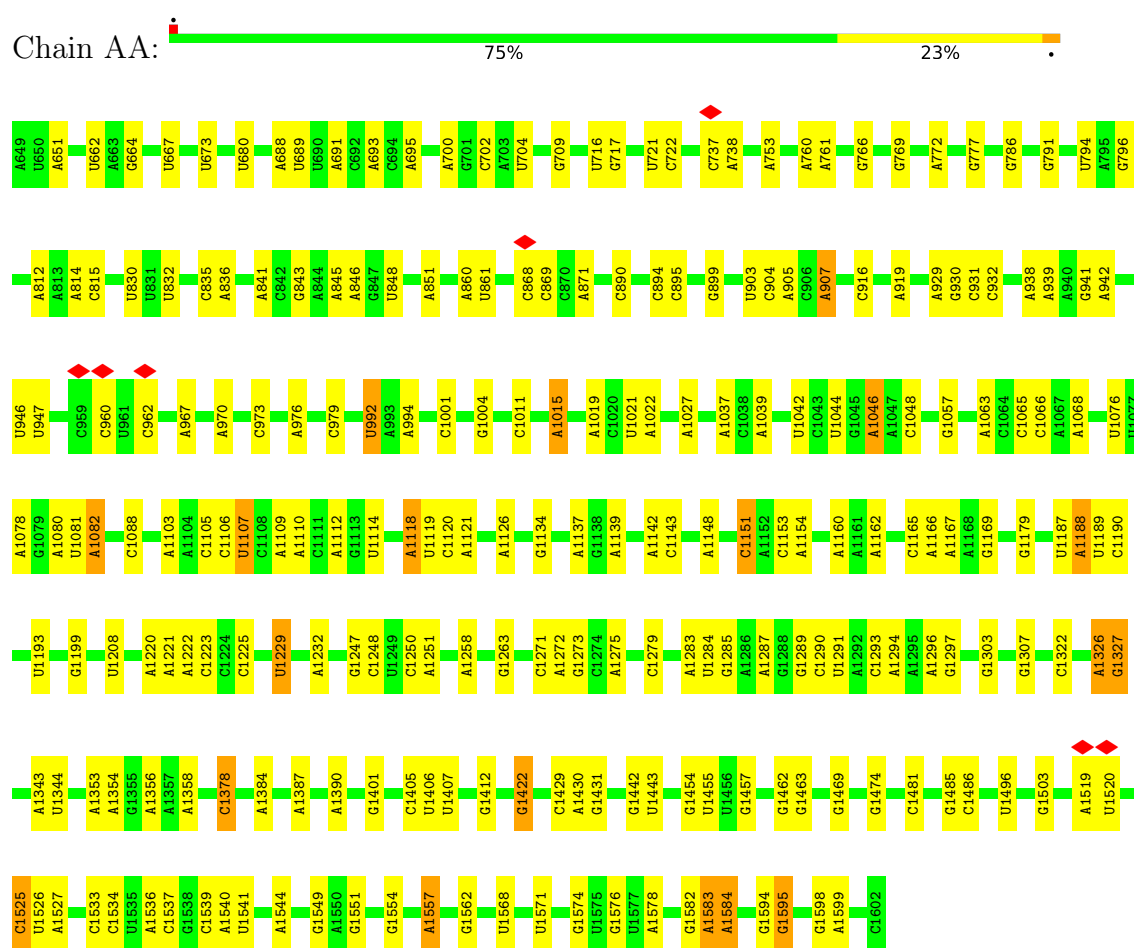
Continued from previous page...

Mol	Chain	Residues	Atoms		AltConf
101	i	46	Total 46	O 46	0
101	j	22	Total 22	O 22	0
101	k	5	Total 5	O 5	0
101	l	5	Total 5	O 5	0
101	m	8	Total 8	O 8	0
101	o	34	Total 34	O 34	0
101	p	8	Total 8	O 8	0
101	r	49	Total 49	O 49	0
101	s	51	Total 51	O 51	0
101	H	7	Total 7	O 7	0

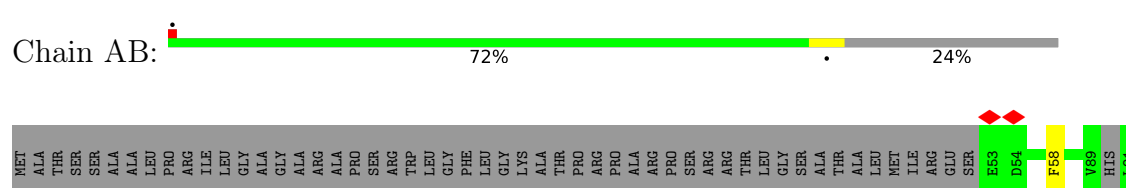
3 Residue-property plots

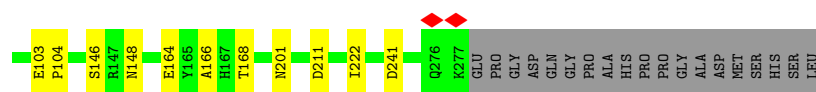
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: 12S mitochondrial rRNA

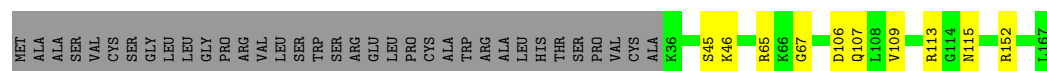
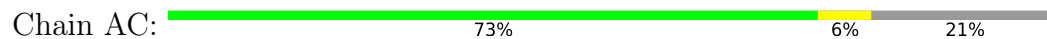


• Molecule 2: 28S ribosomal protein S2, mitochondrial

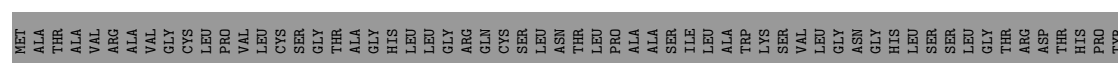
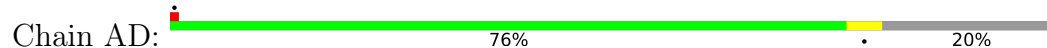




- Molecule 3: 28S ribosomal protein S24, mitochondrial



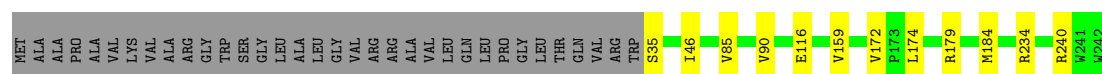
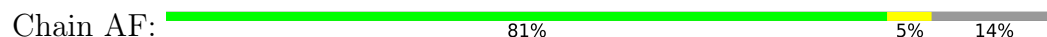
- Molecule 4: 28S ribosomal protein S5, mitochondrial



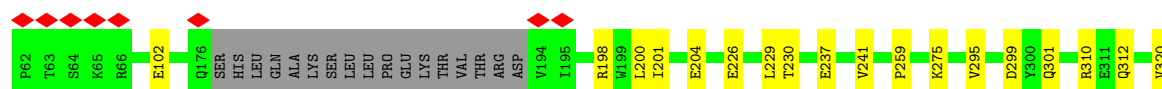
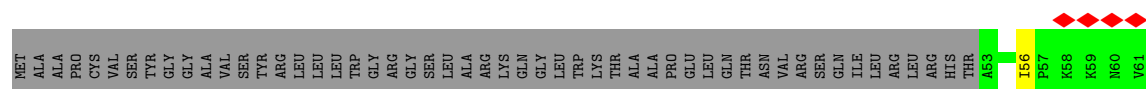
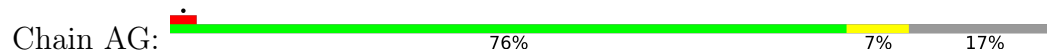
- Molecule 5: 28S ribosomal protein S6, mitochondrial



- Molecule 6: 28S ribosomal protein S7, mitochondrial



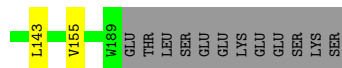
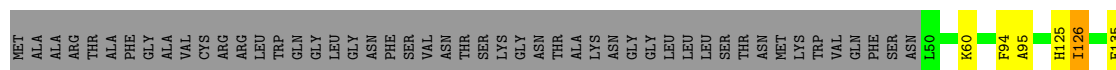
- Molecule 7: 28S ribosomal protein S9, mitochondrial





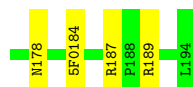
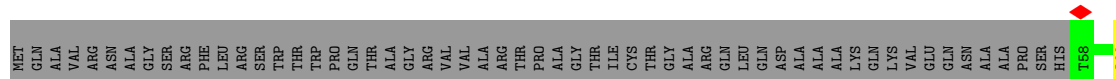
- Molecule 8: 28S ribosomal protein S10, mitochondrial

Chain AH: 66% 30%



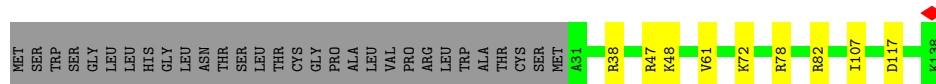
- Molecule 9: 28S ribosomal protein S11, mitochondrial

Chain AI: 68% 29%



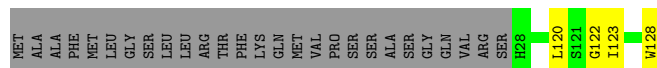
- Molecule 10: 28S ribosomal protein S12, mitochondrial

Chain AJ: 72% 7% 22%



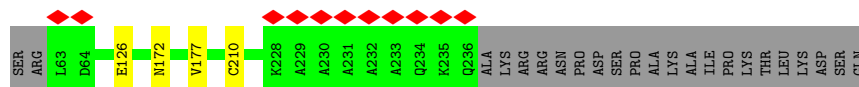
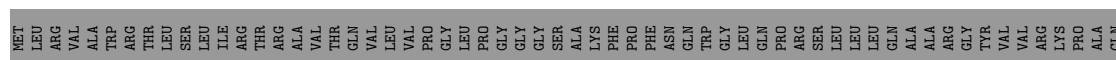
- Molecule 11: 28S ribosomal protein S14, mitochondrial

Chain AK: 76% 21%



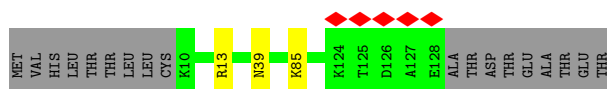
- Molecule 12: 28S ribosomal protein S15, mitochondrial

Chain AL: 66% 32%



- Molecule 13: 28S ribosomal protein S16, mitochondrial

Chain AM:  85% 13%



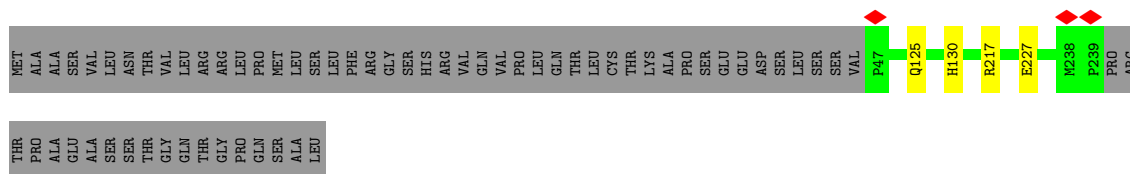
- Molecule 14: 28S ribosomal protein S17, mitochondrial

Chain AN:  79% 5% 15%



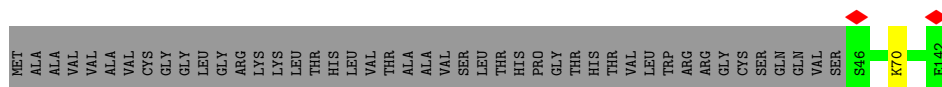
- Molecule 15: 28S ribosomal protein S18b, mitochondrial

Chain AO:  73% 25%

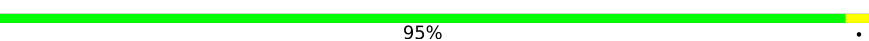


- Molecule 16: 28S ribosomal protein S18c, mitochondrial

Chain AP:  68% 32%



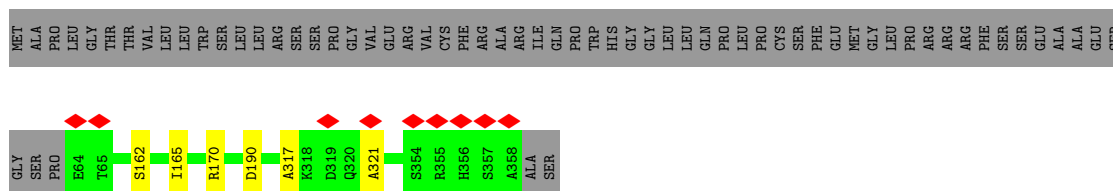
- Molecule 17: MRPS21 isoform 1

Chain AQ:  95%

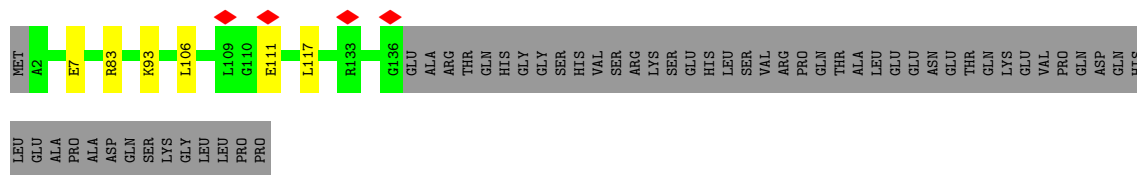


- Molecule 18: 28S ribosomal protein S22, mitochondrial

Chain AR:  80% 18%



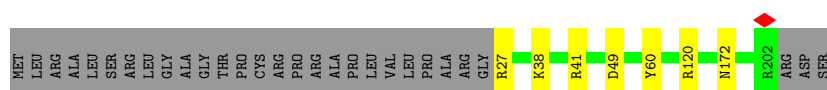
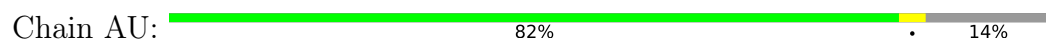
- Molecule 19: 28S ribosomal protein S23, mitochondrial



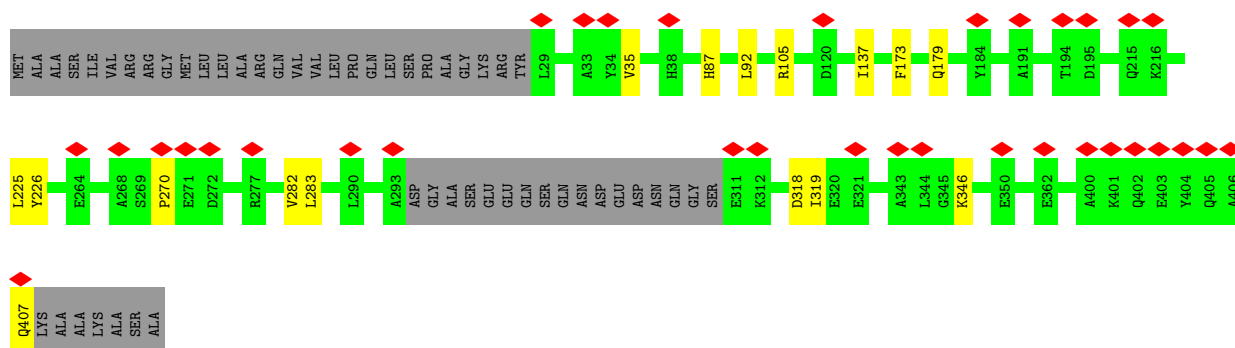
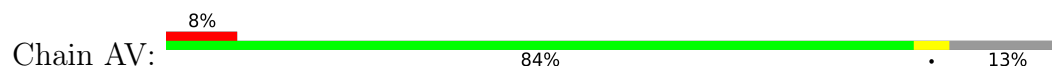
- Molecule 20: 28S ribosomal protein S25, mitochondrial



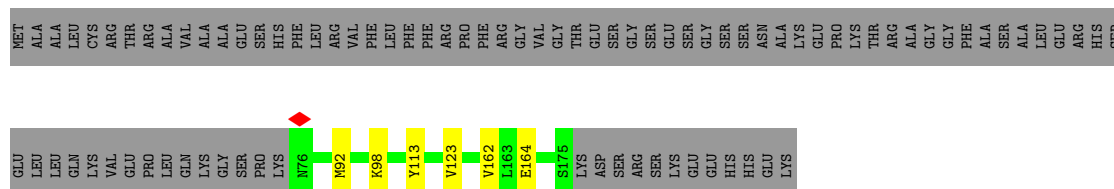
- Molecule 21: 28S ribosomal protein S26, mitochondrial



- Molecule 22: 28S ribosomal protein S27, mitochondrial



- Molecule 23: 28S ribosomal protein S28, mitochondrial

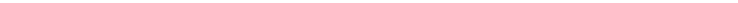


- Molecule 24: 28S ribosomal protein S29, mitochondrial

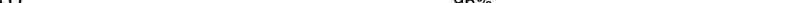
V141	L151	L152	L183	D167	E193	H187	M190	L244	L247	T265	T272	K279	T282	M297	E379	L383	S394	N395	S399	R393	L398																			
Met	Met	Lys	Ile	Thr	Leu	Ser	Arg	His	Lys	Leu	Asp	Pro	Gly	Phe	Leu	His	Met	Gly	Thr	Gln	Ala	Arg	Gln	Ser	Ile	Ala	Ala	His	Leu	Asp	Asn	Gln	Val	Pro	Val	Glu	Ser	Pro	Arg	Arg

- Chain AY: 36% 62%

Amino Acid	Percentage
THR	1.61
ASP	1.61
ASP	1.61
LEU	1.61
LYS	1.61
LYS	1.61
R247	1.61
R255	1.61
L256	1.61
F259	1.61
A263	1.61
V264	1.61
T265	1.61
K266	1.61
E267	1.61
A268	1.61
P269	1.61
E270	1.61
T271	1.61
D272	1.61
T273	1.61
N290	1.61
K309	1.61
E318	1.61
R377	1.61
N381	1.61
N395	1.61

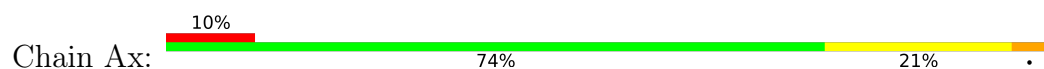
- Chain AZ:  92% 6%

MET
 SER
 S3
 R17
 K46
 Y49
 G99
 A102
 ALA
 LYS
 ARG
 LYS

- Chain A0:  96%

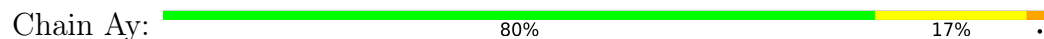
- Chain A1: 

MET	ALA	ALA	ALA	ALA	LEU	PRO	ALA	ALA	TRP	TRP	SER	LEU	GLN	SER	ARG	ALA	ARG	THR	LEU	ALA	PHE	SER	THR	THR	VAL	VAL	THR	PRO	PRO	PRO	PRO	GLU	ARG	THR	THR	PRO	GLY	ASN	E45	R46	P47	R66	P96	L103	H113	L114	D128	F129
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	------	------	------	------	------

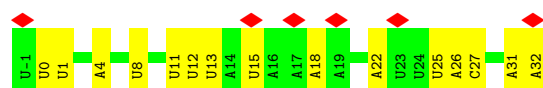




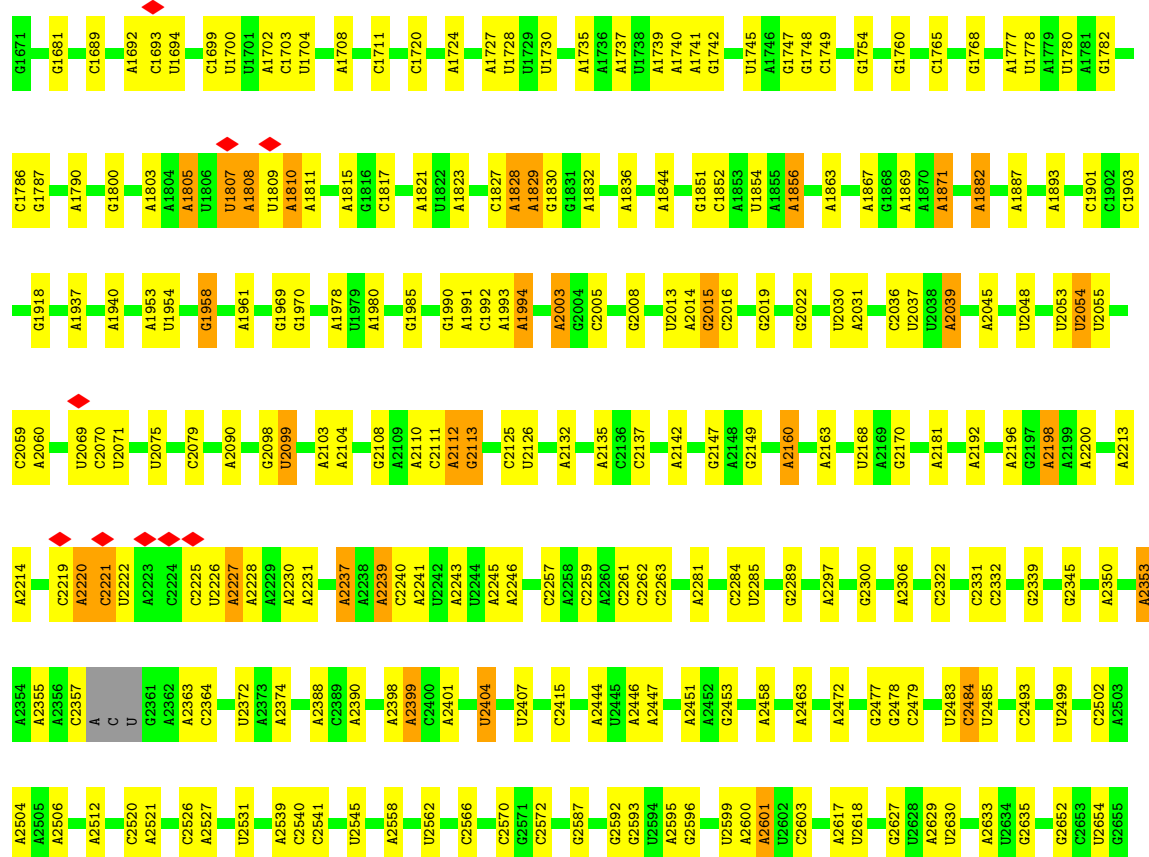
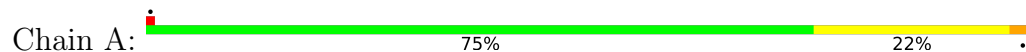
• Molecule 34: E/E-tRNA

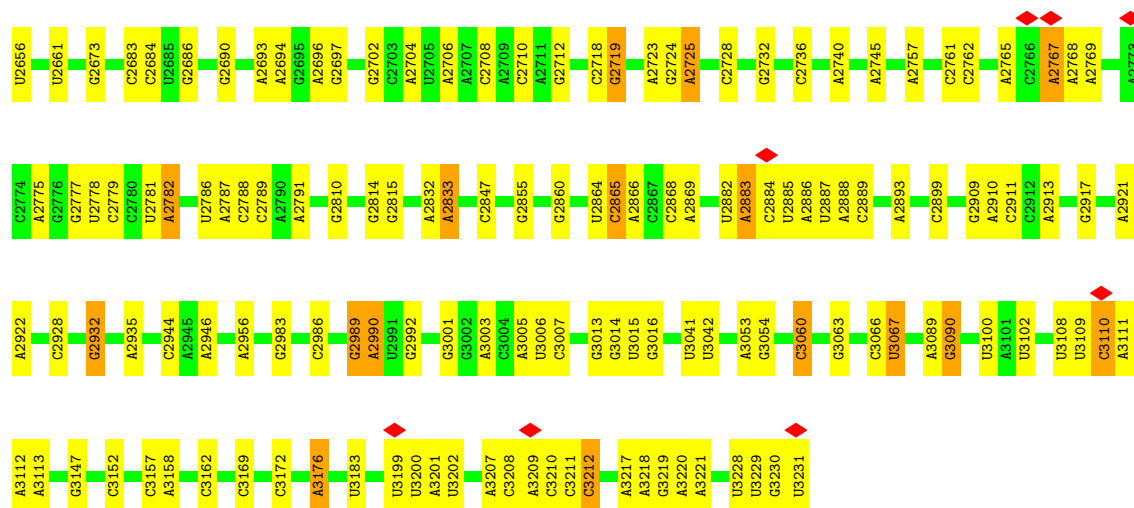


• Molecule 35: mRNA

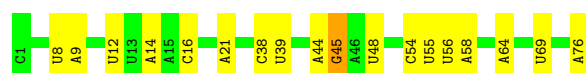


• Molecule 36: 16S mitochondrial rRNA

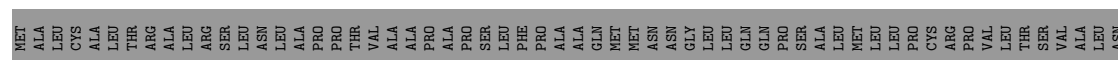
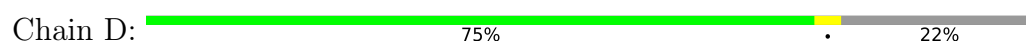




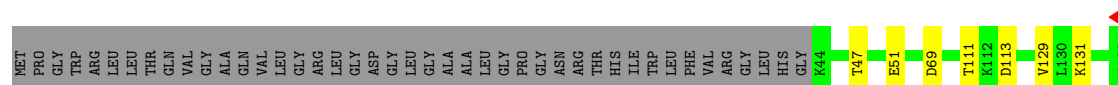
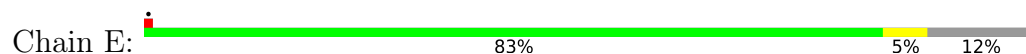
- Molecule 37: mitochondrial tRNAVal



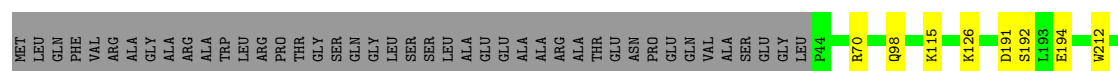
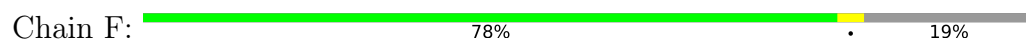
- Molecule 38: 39S ribosomal protein L2, mitochondrial

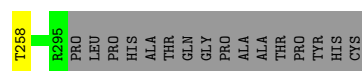


- Molecule 39: 39S ribosomal protein L3, mitochondrial

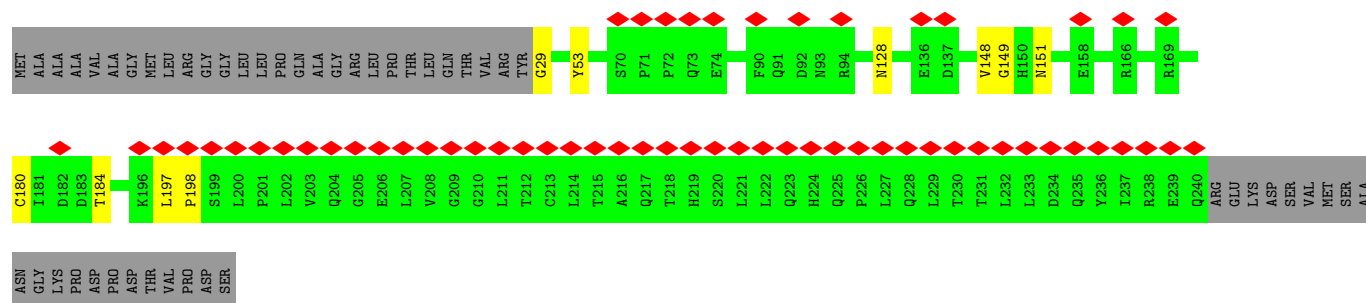
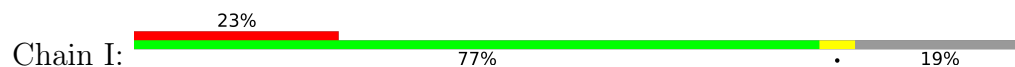


- Molecule 40: 39S ribosomal protein L4, mitochondrial

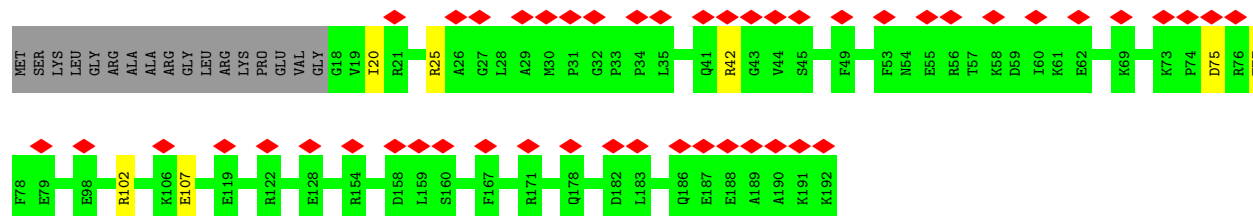
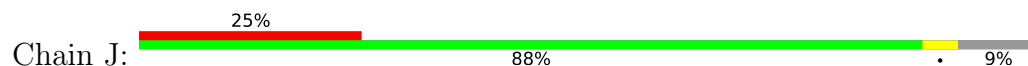




- Molecule 41: 39S ribosomal protein L10, mitochondrial



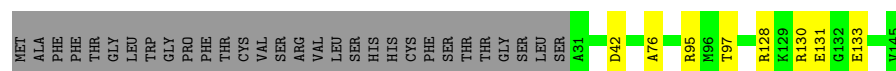
- Molecule 42: 39S ribosomal protein L11, mitochondrial



- Molecule 43: 39S ribosomal protein L13, mitochondrial



- Molecule 44: 39S ribosomal protein L14, mitochondrial

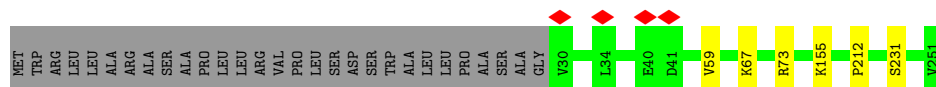


- Molecule 45: 39S ribosomal protein L15, mitochondrial



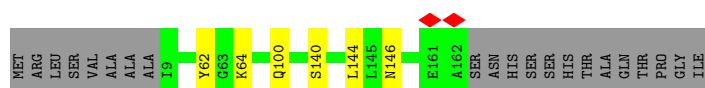
- Molecule 46: 39S ribosomal protein L16, mitochondrial

Chain N:  86% 12%



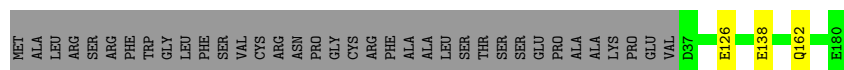
- Molecule 47: 39S ribosomal protein L17, mitochondrial

Chain O:  85% 12%



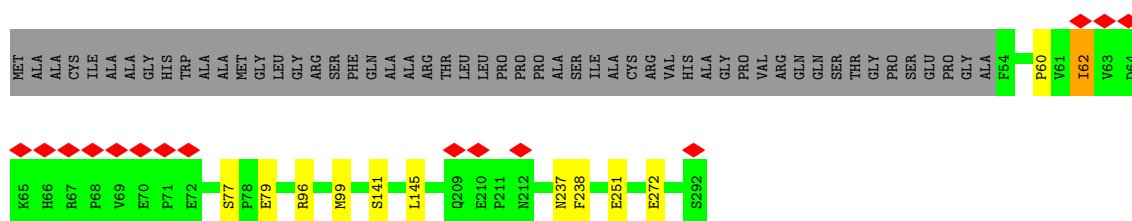
- Molecule 48: 39S ribosomal protein L18, mitochondrial

Chain P:  78% 20%

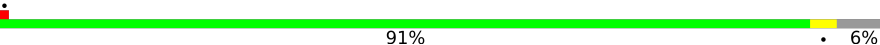


- Molecule 49: 39S ribosomal protein L19, mitochondrial

Chain Q:  5% 78% 18%



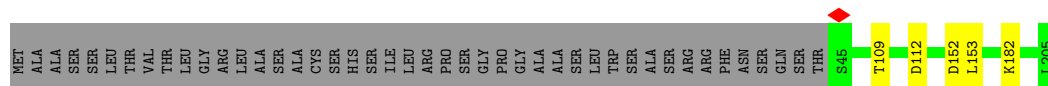
- Molecule 50: 39S ribosomal protein L20, mitochondrial

Chain R:  91% 6%



- Molecule 51: 39S ribosomal protein L21, mitochondrial

Chain S:  76% 21%



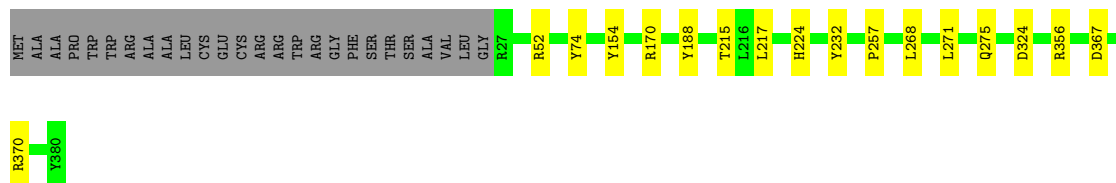
- Molecule 52: 39S ribosomal protein L22, mitochondrial

Chain T:  80% 19%

- Chain 5: 91% 7%
- MET ALA LEU LEU SER SER GLY PRO ALA ARG ARG ALA LEU LEU ALA GLY SER GLY GLN LEU LEU LEU GLY PHE GLY ALA PRO ARG ARG ARG A30 E56 Y105 R110 R144 R173 Y176 C177 P178 I262 D295 D377 R395 K413

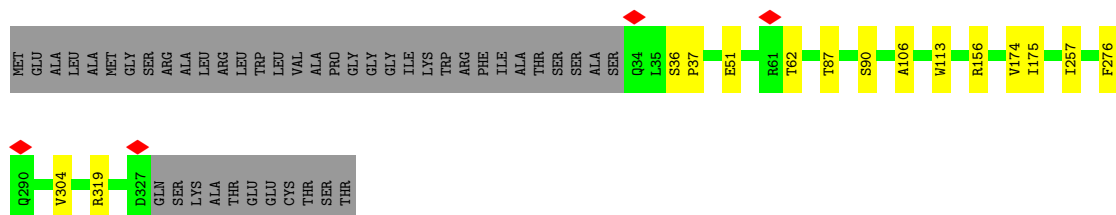
- Molecule 65: 39S ribosomal protein L38, mitochondrial

Chain 6:  89% 7%



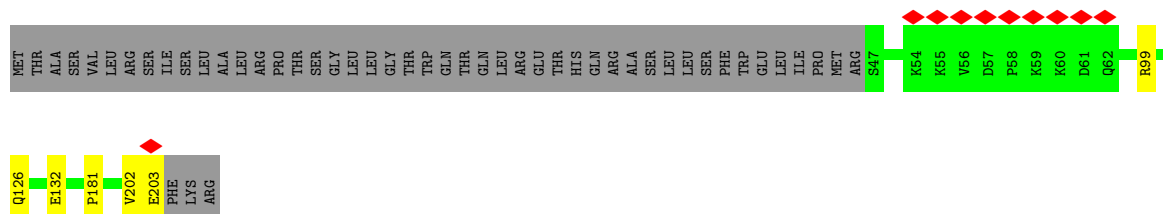
- Molecule 66: 39S ribosomal protein L39, mitochondrial

Chain 7:  83% 13%



- Molecule 67: 39S ribosomal protein L40, mitochondrial

Chain 8:  5% 73% 24%



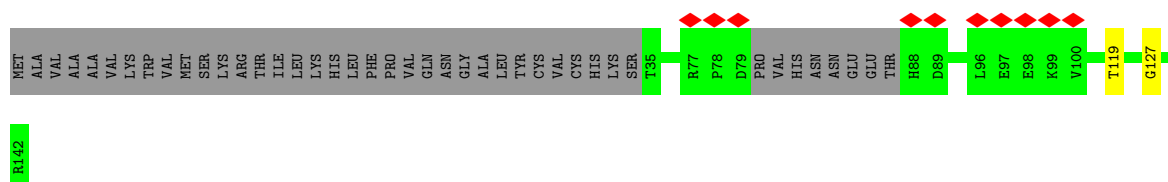
- Molecule 68: 39S ribosomal protein L41, mitochondrial

Chain 9:  88% 9%



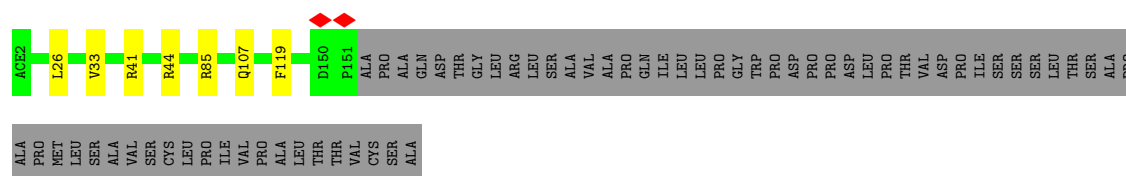
- Molecule 69: 39S ribosomal protein L42, mitochondrial

Chain a:  7% 69% 30%



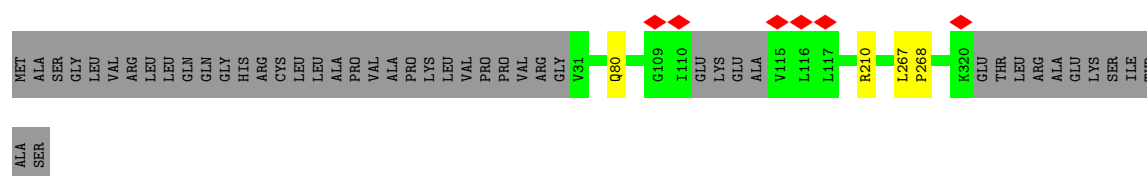
- Molecule 70: Large ribosomal subunit protein mL43

Chain b:  67% 30%



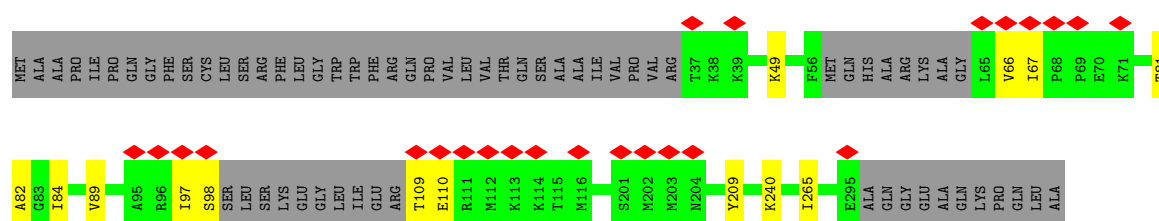
- Molecule 71: 39S ribosomal protein L44, mitochondrial

Chain c:  85% 14%



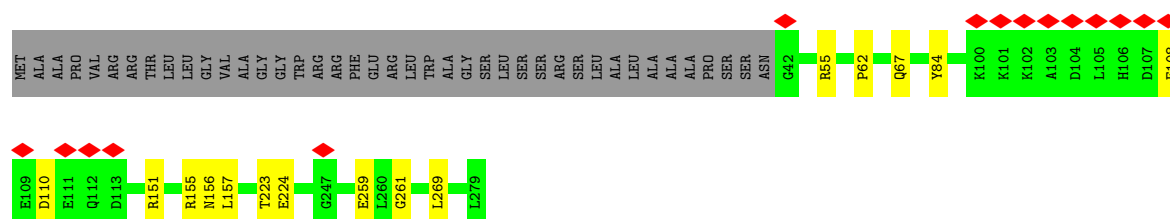
- Molecule 72: 39S ribosomal protein L45, mitochondrial

Chain d:  8% 74% 5% 21%



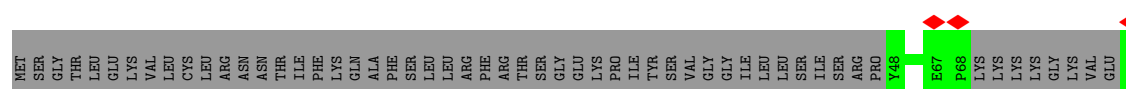
- Molecule 73: 39S ribosomal protein L46, mitochondrial

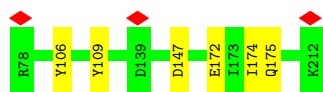
Chain e:  5% 80% 5% 15%



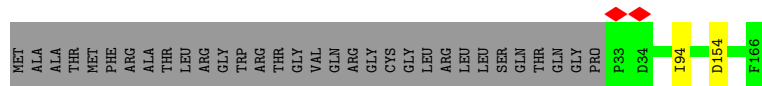
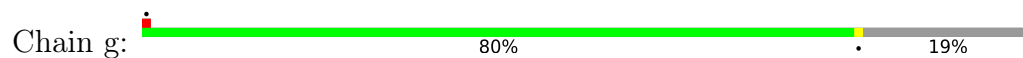
- Molecule 74: 39S ribosomal protein L48, mitochondrial

Chain f:  71% 26%

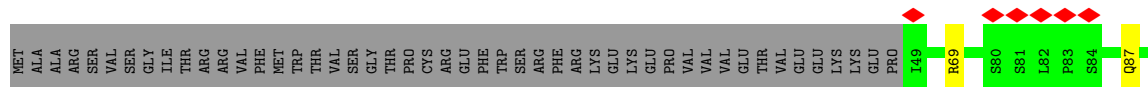




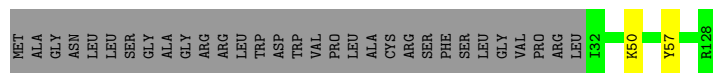
- Molecule 75: 39S ribosomal protein L49, mitochondrial



- Molecule 76: 39S ribosomal protein L50, mitochondrial



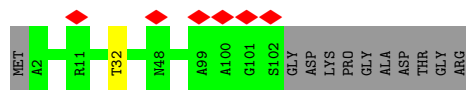
- Molecule 77: 39S ribosomal protein L51, mitochondrial



- Molecule 78: 39S ribosomal protein L52, mitochondrial



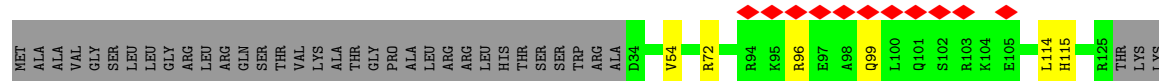
- Molecule 79: 39S ribosomal protein L53, mitochondrial



- Molecule 80: 39S ribosomal protein L54, mitochondrial



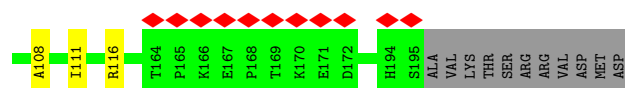
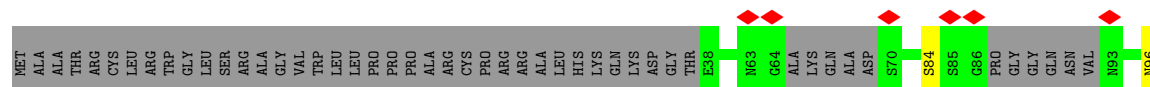
- Molecule 81: 39S ribosomal protein L55, mitochondrial



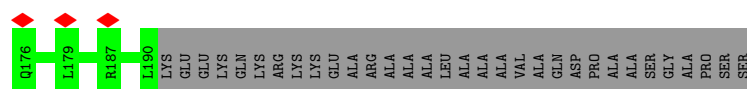
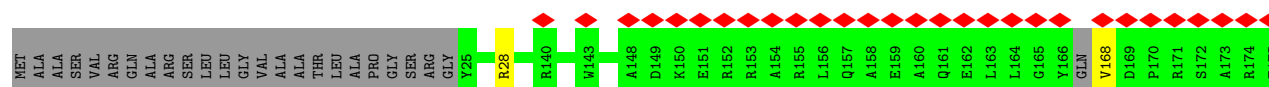
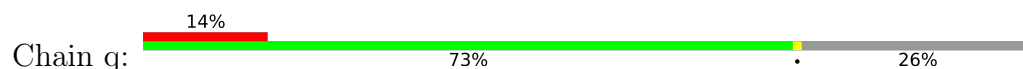
- Molecule 82: Ribosomal protein 63, mitochondrial



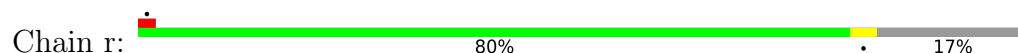
- Molecule 83: Peptidyl-tRNA hydrolase ICT1, mitochondrial

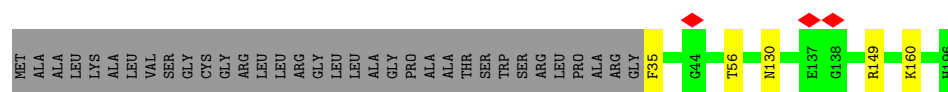


- Molecule 84: Growth arrest and DNA damage-inducible proteins-interacting protein 1



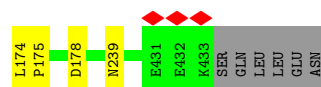
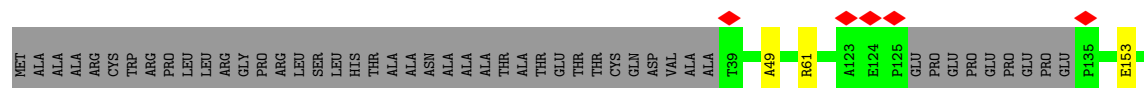
- Molecule 85: 39S ribosomal protein S18a, mitochondrial





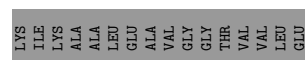
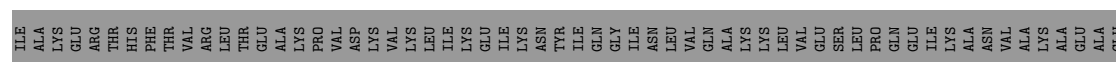
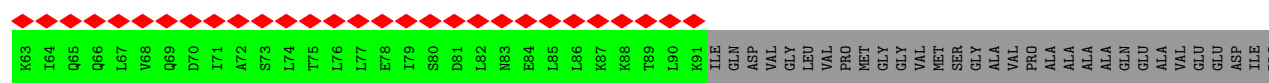
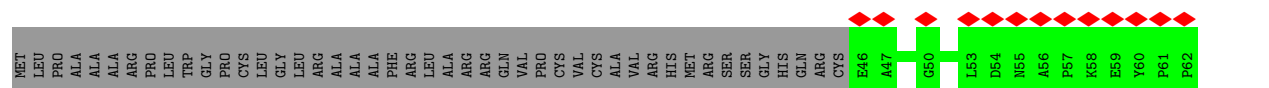
- Molecule 86: 39S ribosomal protein S30, mitochondrial

Chain s: 86% 12%



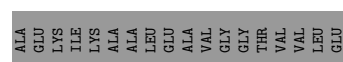
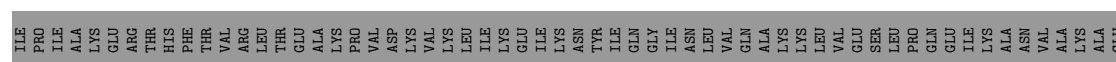
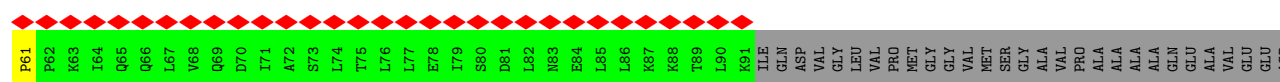
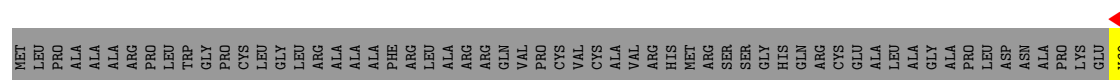
- Molecule 87: 39S ribosomal protein L12, mitochondrial

Chain t: 21% 23% 77%



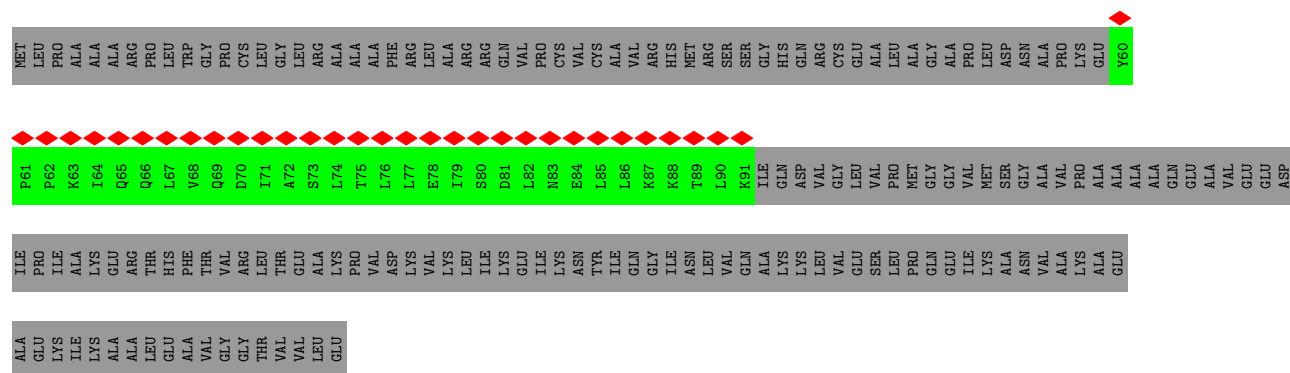
- Molecule 87: 39S ribosomal protein L12, mitochondrial

Chain u: 16% 15% 84%



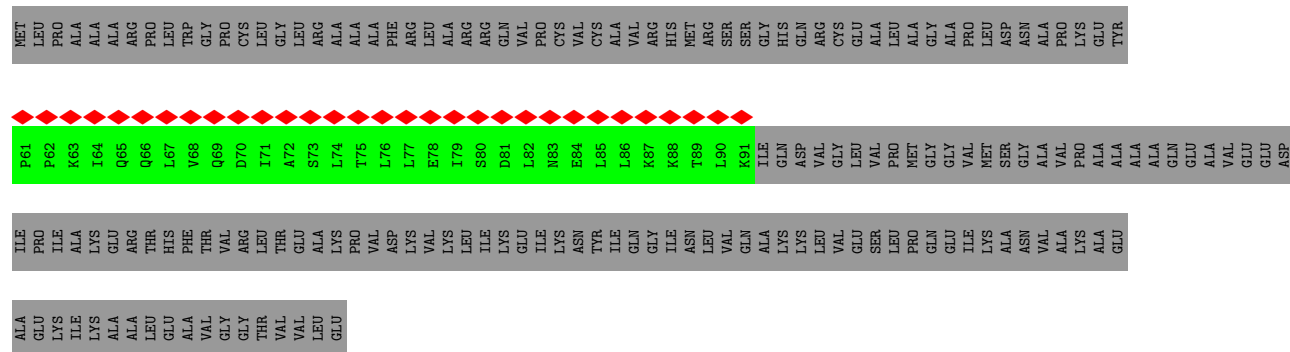
- Molecule 87: 39S ribosomal protein L12, mitochondrial

Chain v: 16% 16% 84%



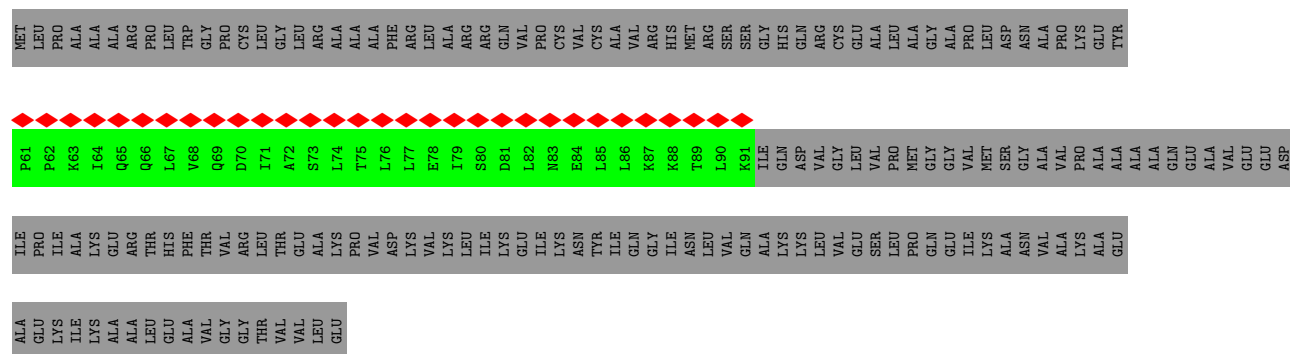
• Molecule 87: 39S ribosomal protein L12, mitochondrial

Chain w: 



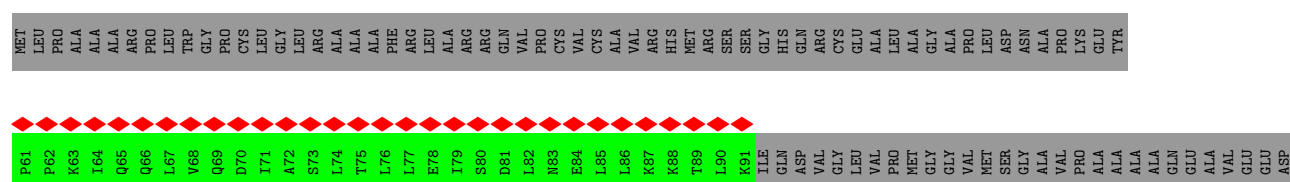
• Molecule 87: 39S ribosomal protein L12, mitochondrial

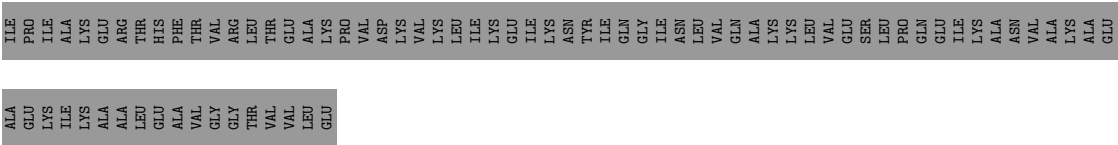
Chain x: 



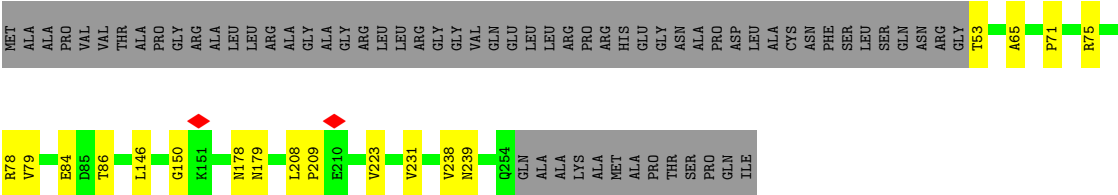
• Molecule 87: 39S ribosomal protein L12, mitochondrial

Chain y: 

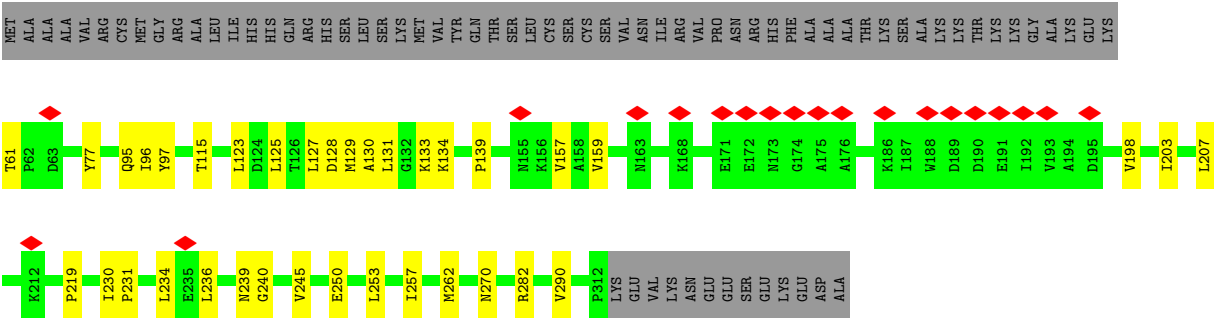




● Molecule 88: 39S ribosomal protein L9, mitochondrial



● Molecule 89: 39S ribosomal protein L1, mitochondrial



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	82522	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 ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	61.722	Depositor
Minimum map value	-28.634	Depositor
Average map value	0.003	Depositor
Map value standard deviation	1.267	Depositor
Recommended contour level	2.8	Depositor
Map size ($\text{\AA}$)	448.19998, 448.19998, 448.19998	wwPDB
Map dimensions	540, 540, 540	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.83, 0.83, 0.83	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: GDP, SPD, 2MG, SAC, ATP, SPM, PUT, B8T, NAD, PSU, OMU, 1MA, ZN, ACE, K, FES, MA6, OMG, AYA, 5MC, MG, 5MU, 5F0

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	AA	0.09	0/22537	0.18	0/35085
2	AB	0.08	0/1859	0.23	0/2513
3	AC	0.11	0/1113	0.26	0/1505
4	AD	0.09	0/2783	0.24	0/3724
5	AE	0.08	0/989	0.23	0/1335
6	AF	0.09	0/1767	0.23	0/2373
7	AG	0.09	0/2746	0.24	0/3681
8	AH	0.13	0/1178	0.28	0/1598
9	AI	0.08	0/1030	0.23	0/1386
10	AJ	0.10	0/855	0.27	0/1148
11	AK	0.10	0/880	0.21	0/1182
12	AL	0.08	0/1477	0.21	0/1974
13	AM	0.09	0/963	0.25	0/1295
14	AN	0.09	0/886	0.26	0/1199
15	AO	0.09	0/1648	0.23	0/2243
16	AP	0.09	0/798	0.23	0/1070
17	AQ	0.08	0/748	0.23	0/994
18	AR	0.07	0/2456	0.20	0/3317
19	AS	0.08	0/1138	0.19	0/1533
20	AT	0.08	0/1402	0.23	0/1883
21	AU	0.07	0/1510	0.18	0/2025
22	AV	0.06	0/3030	0.19	0/4093
23	AW	0.08	0/801	0.23	0/1079
24	AX	0.10	0/2921	0.25	0/3954
25	AY	0.10	0/1280	0.22	0/1725
26	AZ	0.09	0/857	0.23	0/1141
27	A0	0.08	0/1834	0.23	0/2484
28	A1	0.10	0/2313	0.25	0/3129
29	A2	0.10	0/941	0.27	0/1257
30	A3	0.09	0/636	0.28	0/839
31	A4	0.09	0/4877	0.23	0/6598

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	Aw	0.07	0/1603	0.17	0/2488
33	Ax	0.08	0/1655	0.19	0/2569
34	Ay	0.08	0/1656	0.18	0/2571
35	Az	0.10	0/803	0.22	0/1244
36	A	0.09	0/36876	0.21	0/57402
37	B	0.09	0/1627	0.17	0/2527
38	D	0.08	0/1896	0.26	0/2549
39	E	0.10	0/2475	0.27	0/3355
40	F	0.10	0/2090	0.28	0/2842
41	I	0.10	0/1731	0.27	0/2345
42	J	0.09	0/1348	0.23	0/1813
43	K	0.09	0/1490	0.26	0/2021
44	L	0.09	0/905	0.27	0/1218
45	M	0.10	0/2368	0.29	0/3195
46	N	0.09	0/1833	0.27	0/2468
47	O	0.08	0/1283	0.25	0/1727
48	P	0.11	0/1199	0.26	0/1623
49	Q	0.09	0/2039	0.25	0/2750
50	R	0.09	0/1175	0.23	0/1572
51	S	0.10	0/1320	0.29	0/1789
52	T	0.10	0/1403	0.27	0/1886
53	U	0.10	0/1274	0.25	0/1723
54	V	0.11	0/1721	0.28	0/2333
55	W	0.10	0/926	0.26	0/1244
56	X	0.10	0/2099	0.23	0/2837
57	Y	0.09	0/1593	0.22	0/2136
58	Z	0.09	0/1021	0.27	0/1378
59	0	0.09	0/913	0.24	0/1224
60	1	0.09	0/469	0.27	0/621
61	2	0.10	0/383	0.29	0/507
62	3	0.10	0/853	0.27	0/1136
63	4	0.08	0/350	0.22	0/461
64	5	0.09	0/3305	0.26	0/4502
65	6	0.11	0/3043	0.26	0/4140
66	7	0.10	0/2447	0.26	0/3310
67	8	0.09	0/1354	0.25	0/1819
68	9	0.09	0/1025	0.25	0/1379
69	a	0.10	0/866	0.29	0/1174
70	b	0.10	0/1219	0.26	0/1651
71	c	0.09	0/2347	0.22	0/3171
72	d	0.09	0/2039	0.25	0/2759
73	e	0.09	0/1970	0.22	0/2658
74	f	0.09	0/1273	0.21	0/1716

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
75	g	0.10	0/1151	0.26	0/1569
76	h	0.08	0/918	0.21	0/1249
77	i	0.10	0/850	0.26	0/1135
78	j	0.11	0/760	0.25	0/1023
79	k	0.07	0/777	0.21	0/1048
80	l	0.08	0/707	0.25	0/960
81	m	0.08	0/805	0.22	0/1081
82	o	0.11	0/819	0.28	0/1097
83	p	0.08	0/1223	0.23	0/1641
84	q	0.08	0/1422	0.22	0/1916
85	r	0.10	0/1362	0.28	0/1846
86	s	0.09	0/3239	0.26	0/4400
87	t	0.06	0/358	0.19	0/486
87	u	0.08	0/259	0.20	0/350
87	v	0.07	0/259	0.16	0/350
87	w	0.08	0/246	0.19	0/331
87	x	0.07	0/246	0.17	0/331
87	y	0.07	0/246	0.17	0/331
88	H	0.10	0/1698	0.26	0/2292
89	z	0.11	0/2067	0.32	0/2793
All	All	0.09	0/190930	0.23	0/271424

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	AA	20260	10304	10285	109	0
2	AB	1818	1809	1807	11	0
3	AC	1083	1089	1088	8	0
4	AD	2731	2805	2804	13	0
5	AE	972	1000	1000	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	AF	1725	1771	1769	12	0
7	AG	2688	2689	2687	18	0
8	AH	1152	1187	1183	7	0
9	AI	1020	1061	1053	4	0
10	AJ	839	889	887	10	0
11	AK	862	886	885	3	0
12	AL	1453	1541	1540	3	0
13	AM	942	966	965	3	0
14	AN	868	929	928	4	0
15	AO	1592	1561	1557	3	0
16	AP	781	807	806	1	0
17	AQ	744	758	758	4	0
18	AR	2409	2430	2428	4	0
19	AS	1111	1116	1115	4	0
20	AT	1371	1393	1393	4	0
21	AU	1488	1501	1499	7	0
22	AV	2969	2964	2961	11	0
23	AW	789	804	802	4	0
24	AX	2849	2845	2843	21	0
25	AY	1246	1198	1197	6	0
26	AZ	839	859	858	2	0
27	A0	1787	1797	1796	4	0
28	A1	2265	2296	2294	14	0
29	A2	935	971	971	11	0
30	A3	625	701	699	4	0
31	A4	4768	4770	4766	14	0
32	Aw	1434	725	725	0	0
33	Ax	1482	751	752	7	0
34	Ay	1483	752	753	3	0
35	Az	719	360	360	1	0
36	A	33070	16801	16795	169	0
37	B	1524	779	779	3	0
38	D	1859	1921	1920	7	0
39	E	2406	2416	2415	12	0
40	F	2031	2066	2065	6	0
41	I	1695	1789	1785	8	0
42	J	1330	1409	1407	4	0
43	K	1455	1452	1452	5	0
44	L	890	942	941	7	0
45	M	2314	2385	2384	9	0
46	N	1786	1818	1817	6	0
47	O	1259	1295	1294	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
48	P	1173	1166	1165	3	0
49	Q	1990	2031	2031	9	0
50	R	1154	1215	1214	3	0
51	S	1293	1366	1365	5	0
52	T	1369	1412	1410	2	0
53	U	1251	1232	1232	7	0
54	V	1676	1689	1687	12	0
55	W	904	936	935	2	0
56	X	2044	2061	2060	4	0
57	Y	1556	1598	1597	2	0
58	Z	996	1045	1044	4	0
59	0	898	917	916	5	0
60	1	464	513	511	3	0
61	2	377	407	406	1	0
62	3	832	884	883	6	0
63	4	342	362	361	2	0
64	5	3210	3209	3206	7	0
65	6	2948	2844	2841	13	0
66	7	2390	2399	2397	10	0
67	8	1327	1369	1368	5	0
68	9	997	988	987	3	0
69	a	840	812	810	2	0
70	b	1196	1196	1195	6	0
71	c	2299	2322	2320	2	0
72	d	1985	1979	1976	10	0
73	e	1931	1917	1916	10	0
74	f	1252	1271	1269	4	0
75	g	1113	1097	1097	2	0
76	h	895	882	881	2	0
77	i	828	859	857	2	0
78	j	745	747	746	5	0
79	k	774	785	784	1	0
80	l	688	675	674	3	0
81	m	791	760	796	5	0
82	o	798	806	804	3	0
83	p	1205	1224	1223	3	0
84	q	1389	1376	1374	2	0
85	r	1322	1349	1348	5	0
86	s	3155	3143	3139	4	0
87	t	354	378	377	0	0
87	u	257	284	283	1	0
87	v	257	284	283	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
87	w	245	275	275	0	0
87	x	245	275	275	0	0
87	y	245	275	275	0	0
88	H	1661	1736	1734	12	0
89	z	2027	2077	2076	28	0
90	AA	44	26	26	1	0
91	AA	14	30	26	0	0
92	A	30	66	57	0	0
92	AA	10	22	19	0	0
93	A	138	0	0	0	0
93	A3	1	0	0	0	0
93	AA	63	0	0	0	0
93	AB	1	0	0	0	0
93	AX	1	0	0	0	0
93	D	2	0	0	0	0
93	E	1	0	0	0	0
93	g	1	0	0	0	0
94	3	1	0	0	0	0
94	6	1	0	0	0	0
94	A	29	0	0	0	0
94	AA	18	0	0	0	0
94	D	1	0	0	0	0
94	M	1	0	0	0	0
94	W	1	0	0	0	0
94	i	1	0	0	0	0
94	o	1	0	0	0	0
95	0	1	0	0	0	0
95	4	1	0	0	0	0
95	AO	1	0	0	0	0
96	AP	4	0	0	0	0
96	AT	4	0	0	0	0
96	r	4	0	0	0	0
97	AX	31	12	12	0	0
98	AX	28	12	12	0	0
99	A	6	14	12	0	0
100	B	7	11	8	0	0
101	0	28	0	0	2	0
101	1	3	0	0	0	0
101	2	34	0	0	0	0
101	3	44	0	0	2	0
101	4	11	0	0	0	0
101	5	14	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
101	6	100	0	0	4	0
101	7	15	0	0	0	0
101	8	15	0	0	2	0
101	9	17	0	0	0	0
101	A	2955	0	0	87	0
101	A0	1	0	0	0	0
101	A1	35	0	0	0	0
101	A2	32	0	0	0	0
101	A3	50	0	0	1	0
101	A4	9	0	0	0	0
101	AA	2039	0	0	59	0
101	AB	78	0	0	7	0
101	AC	55	0	0	3	0
101	AD	82	0	0	2	0
101	AE	27	0	0	0	0
101	AF	39	0	0	2	0
101	AG	64	0	0	3	0
101	AH	57	0	0	1	0
101	AI	39	0	0	2	0
101	AJ	25	0	0	2	0
101	AK	51	0	0	2	0
101	AL	42	0	0	2	0
101	AM	22	0	0	1	0
101	AN	39	0	0	0	0
101	AO	37	0	0	2	0
101	AP	31	0	0	1	0
101	AQ	77	0	0	1	0
101	AR	11	0	0	1	0
101	AS	25	0	0	0	0
101	AT	44	0	0	0	0
101	AU	8	0	0	2	0
101	AW	16	0	0	0	0
101	AX	56	0	0	4	0
101	AY	18	0	0	1	0
101	AZ	30	0	0	0	0
101	Aw	6	0	0	0	0
101	Ax	7	0	0	0	0
101	Ay	3	0	0	0	0
101	Az	13	0	0	0	0
101	B	66	0	0	0	0
101	D	68	0	0	0	0
101	E	66	0	0	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
101	F	79	0	0	3	0
101	H	7	0	0	1	0
101	I	27	0	0	3	0
101	J	1	0	0	0	0
101	K	57	0	0	2	0
101	L	31	0	0	4	0
101	M	51	0	0	2	0
101	N	58	0	0	1	0
101	O	33	0	0	0	0
101	P	81	0	0	3	0
101	Q	29	0	0	1	0
101	R	63	0	0	2	0
101	S	49	0	0	2	0
101	T	45	0	0	1	0
101	U	24	0	0	0	0
101	V	5	0	0	0	0
101	W	55	0	0	1	0
101	X	11	0	0	0	0
101	Y	25	0	0	0	0
101	Z	32	0	0	0	0
101	a	11	0	0	1	0
101	b	35	0	0	2	0
101	c	18	0	0	0	0
101	d	9	0	0	0	0
101	e	13	0	0	2	0
101	f	19	0	0	0	0
101	g	13	0	0	1	0
101	i	46	0	0	0	0
101	j	22	0	0	2	0
101	k	5	0	0	1	0
101	l	5	0	0	1	0
101	m	8	0	0	0	0
101	o	34	0	0	2	0
101	p	8	0	0	0	0
101	r	49	0	0	3	0
101	s	51	0	0	0	0
All	All	189566	153978	153843	695	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 2.

The worst 5 of 695 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)
1:AA:1046:A:O2'	1:AA:1048:C:OP2	1.97	0.82
74:f:106:TYR:OH	74:f:174:ILE:O	1.97	0.82
1:AA:1327:G:O6	101:AA:1801:HOH:O	1.99	0.80
36:A:3063:G:O2'	36:A:3066:C:OP2	1.98	0.80
36:A:3042:U:OP2	101:A:3501:HOH:O	1.99	0.78

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	AB	220/296 (74%)	218 (99%)	2 (1%)	0	100	100
3	AC	130/167 (78%)	125 (96%)	5 (4%)	0	100	100
4	AD	341/430 (79%)	331 (97%)	10 (3%)	0	100	100
5	AE	120/125 (96%)	119 (99%)	1 (1%)	0	100	100
6	AF	206/242 (85%)	204 (99%)	2 (1%)	0	100	100
7	AG	323/396 (82%)	315 (98%)	8 (2%)	0	100	100
8	AH	138/201 (69%)	137 (99%)	0	1 (1%)	18	27
9	AI	134/194 (69%)	133 (99%)	1 (1%)	0	100	100
10	AJ	106/138 (77%)	105 (99%)	1 (1%)	0	100	100
11	AK	99/128 (77%)	99 (100%)	0	0	100	100
12	AL	172/257 (67%)	170 (99%)	2 (1%)	0	100	100
13	AM	117/137 (85%)	117 (100%)	0	0	100	100
14	AN	108/130 (83%)	106 (98%)	2 (2%)	0	100	100
15	AO	191/258 (74%)	188 (98%)	3 (2%)	0	100	100
16	AP	95/142 (67%)	94 (99%)	1 (1%)	0	100	100
17	AQ	84/87 (97%)	82 (98%)	2 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
18	AR	293/360 (81%)	284 (97%)	9 (3%)	0	100	100
19	AS	133/190 (70%)	133 (100%)	0	0	100	100
20	AT	166/173 (96%)	164 (99%)	2 (1%)	0	100	100
21	AU	174/205 (85%)	174 (100%)	0	0	100	100
22	AV	358/414 (86%)	351 (98%)	7 (2%)	0	100	100
23	AW	98/187 (52%)	95 (97%)	3 (3%)	0	100	100
24	AX	350/398 (88%)	342 (98%)	7 (2%)	1 (0%)	36	50
25	AY	147/395 (37%)	147 (100%)	0	0	100	100
26	AZ	98/106 (92%)	97 (99%)	1 (1%)	0	100	100
27	A0	213/217 (98%)	209 (98%)	4 (2%)	0	100	100
28	A1	277/323 (86%)	273 (99%)	4 (1%)	0	100	100
29	A2	115/118 (98%)	115 (100%)	0	0	100	100
30	A3	68/199 (34%)	67 (98%)	1 (2%)	0	100	100
31	A4	584/689 (85%)	577 (99%)	7 (1%)	0	100	100
38	D	236/305 (77%)	231 (98%)	5 (2%)	0	100	100
39	E	303/348 (87%)	298 (98%)	4 (1%)	1 (0%)	36	50
40	F	250/311 (80%)	245 (98%)	5 (2%)	0	100	100
41	I	210/261 (80%)	200 (95%)	10 (5%)	0	100	100
42	J	173/192 (90%)	173 (100%)	0	0	100	100
43	K	175/178 (98%)	173 (99%)	2 (1%)	0	100	100
44	L	113/145 (78%)	112 (99%)	1 (1%)	0	100	100
45	M	287/296 (97%)	284 (99%)	3 (1%)	0	100	100
46	N	220/251 (88%)	218 (99%)	2 (1%)	0	100	100
47	O	152/175 (87%)	149 (98%)	3 (2%)	0	100	100
48	P	142/180 (79%)	142 (100%)	0	0	100	100
49	Q	237/292 (81%)	235 (99%)	1 (0%)	1 (0%)	30	42
50	R	138/149 (93%)	138 (100%)	0	0	100	100
51	S	159/205 (78%)	156 (98%)	3 (2%)	0	100	100
52	T	164/206 (80%)	164 (100%)	0	0	100	100
53	U	150/153 (98%)	148 (99%)	2 (1%)	0	100	100
54	V	203/216 (94%)	199 (98%)	4 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
55	W	114/148 (77%)	113 (99%)	1 (1%)	0	100	100
56	X	242/256 (94%)	241 (100%)	1 (0%)	0	100	100
57	Y	179/250 (72%)	176 (98%)	3 (2%)	0	100	100
58	Z	120/161 (74%)	119 (99%)	1 (1%)	0	100	100
59	0	108/188 (57%)	108 (100%)	0	0	100	100
60	1	54/65 (83%)	54 (100%)	0	0	100	100
61	2	44/92 (48%)	43 (98%)	1 (2%)	0	100	100
62	3	93/188 (50%)	92 (99%)	1 (1%)	0	100	100
63	4	36/103 (35%)	36 (100%)	0	0	100	100
64	5	392/423 (93%)	384 (98%)	8 (2%)	0	100	100
65	6	352/380 (93%)	344 (98%)	8 (2%)	0	100	100
66	7	292/338 (86%)	284 (97%)	8 (3%)	0	100	100
67	8	155/206 (75%)	153 (99%)	2 (1%)	0	100	100
68	9	122/137 (89%)	121 (99%)	1 (1%)	0	100	100
69	a	96/142 (68%)	95 (99%)	1 (1%)	0	100	100
70	b	149/215 (69%)	147 (99%)	2 (1%)	0	100	100
71	c	282/332 (85%)	279 (99%)	3 (1%)	0	100	100
72	d	235/306 (77%)	231 (98%)	4 (2%)	0	100	100
73	e	236/279 (85%)	230 (98%)	6 (2%)	0	100	100
74	f	153/212 (72%)	151 (99%)	2 (1%)	0	100	100
75	g	132/166 (80%)	130 (98%)	2 (2%)	0	100	100
76	h	108/158 (68%)	108 (100%)	0	0	100	100
77	i	95/128 (74%)	93 (98%)	2 (2%)	0	100	100
78	j	92/123 (75%)	90 (98%)	2 (2%)	0	100	100
79	k	99/112 (88%)	99 (100%)	0	0	100	100
80	l	80/138 (58%)	79 (99%)	1 (1%)	0	100	100
81	m	90/128 (70%)	89 (99%)	1 (1%)	0	100	100
82	o	92/102 (90%)	91 (99%)	1 (1%)	0	100	100
83	p	141/206 (68%)	140 (99%)	1 (1%)	0	100	100
84	q	161/222 (72%)	160 (99%)	1 (1%)	0	100	100
85	r	160/196 (82%)	159 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
86	s	382/439 (87%)	375 (98%)	7 (2%)	0	100	100
87	t	44/198 (22%)	44 (100%)	0	0	100	100
87	u	30/198 (15%)	30 (100%)	0	0	100	100
87	v	30/198 (15%)	30 (100%)	0	0	100	100
87	w	29/198 (15%)	29 (100%)	0	0	100	100
87	x	29/198 (15%)	29 (100%)	0	0	100	100
87	y	29/198 (15%)	29 (100%)	0	0	100	100
88	H	200/267 (75%)	194 (97%)	6 (3%)	0	100	100
89	z	250/325 (77%)	237 (95%)	12 (5%)	1 (0%)	30	42
All	All	14697/19484 (75%)	14472 (98%)	220 (2%)	5 (0%)	100	100

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	AH	126	ILE
89	z	128	ASP
24	AX	81	HIS
39	E	150	LYS
49	Q	62	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	
2	AB	197/249 (79%)	197 (100%)	0	100	100
3	AC	115/143 (80%)	115 (100%)	0	100	100
4	AD	286/357 (80%)	286 (100%)	0	100	100
5	AE	104/107 (97%)	104 (100%)	0	100	100
6	AF	185/209 (88%)	185 (100%)	0	100	100
7	AG	285/342 (83%)	285 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	AH	130/180 (72%)	130 (100%)	0	100	100
9	AI	104/146 (71%)	104 (100%)	0	100	100
10	AJ	93/118 (79%)	93 (100%)	0	100	100
11	AK	91/113 (80%)	91 (100%)	0	100	100
12	AL	158/226 (70%)	158 (100%)	0	100	100
13	AM	97/113 (86%)	97 (100%)	0	100	100
14	AN	96/115 (84%)	96 (100%)	0	100	100
15	AO	174/230 (76%)	174 (100%)	0	100	100
16	AP	88/123 (72%)	88 (100%)	0	100	100
17	AQ	78/79 (99%)	78 (100%)	0	100	100
18	AR	264/318 (83%)	264 (100%)	0	100	100
19	AS	116/164 (71%)	116 (100%)	0	100	100
20	AT	153/157 (98%)	153 (100%)	0	100	100
21	AU	152/174 (87%)	152 (100%)	0	100	100
22	AV	325/364 (89%)	325 (100%)	0	100	100
23	AW	87/158 (55%)	87 (100%)	0	100	100
24	AX	311/351 (89%)	311 (100%)	0	100	100
25	AY	137/357 (38%)	137 (100%)	0	100	100
26	AZ	90/95 (95%)	90 (100%)	0	100	100
27	A0	188/189 (100%)	188 (100%)	0	100	100
28	A1	257/291 (88%)	257 (100%)	0	100	100
29	A2	100/101 (99%)	100 (100%)	0	100	100
30	A3	65/166 (39%)	65 (100%)	0	100	100
31	A4	526/609 (86%)	526 (100%)	0	100	100
38	D	192/245 (78%)	192 (100%)	0	100	100
39	E	260/290 (90%)	260 (100%)	0	100	100
40	F	219/262 (84%)	219 (100%)	0	100	100
41	I	194/232 (84%)	194 (100%)	0	100	100
42	J	138/150 (92%)	138 (100%)	0	100	100
43	K	154/155 (99%)	154 (100%)	0	100	100
44	L	98/124 (79%)	98 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
45	M	245/249 (98%)	245 (100%)	0	100	100
46	N	189/211 (90%)	189 (100%)	0	100	100
47	O	134/150 (89%)	134 (100%)	0	100	100
48	P	126/155 (81%)	126 (100%)	0	100	100
49	Q	221/256 (86%)	221 (100%)	0	100	100
50	R	118/126 (94%)	118 (100%)	0	100	100
51	S	146/180 (81%)	146 (100%)	0	100	100
52	T	146/176 (83%)	146 (100%)	0	100	100
53	U	134/135 (99%)	134 (100%)	0	100	100
54	V	183/191 (96%)	183 (100%)	0	100	100
55	W	94/119 (79%)	94 (100%)	0	100	100
56	X	220/229 (96%)	220 (100%)	0	100	100
57	Y	163/223 (73%)	163 (100%)	0	100	100
58	Z	113/147 (77%)	113 (100%)	0	100	100
59	0	99/164 (60%)	99 (100%)	0	100	100
60	1	53/60 (88%)	53 (100%)	0	100	100
61	2	40/72 (56%)	40 (100%)	0	100	100
62	3	88/166 (53%)	88 (100%)	0	100	100
63	4	37/89 (42%)	37 (100%)	0	100	100
64	5	353/368 (96%)	353 (100%)	0	100	100
65	6	313/332 (94%)	313 (100%)	0	100	100
66	7	270/303 (89%)	270 (100%)	0	100	100
67	8	146/190 (77%)	146 (100%)	0	100	100
68	9	104/112 (93%)	104 (100%)	0	100	100
69	a	96/133 (72%)	96 (100%)	0	100	100
70	b	132/185 (71%)	132 (100%)	0	100	100
71	c	251/288 (87%)	251 (100%)	0	100	100
72	d	223/274 (81%)	223 (100%)	0	100	100
73	e	207/236 (88%)	207 (100%)	0	100	100
74	f	139/188 (74%)	139 (100%)	0	100	100
75	g	124/148 (84%)	124 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
76	h	104/148 (70%)	104 (100%)	0	100	100
77	i	86/110 (78%)	86 (100%)	0	100	100
78	j	74/97 (76%)	74 (100%)	0	100	100
79	k	83/90 (92%)	83 (100%)	0	100	100
80	l	76/116 (66%)	76 (100%)	0	100	100
81	m	85/113 (75%)	85 (100%)	0	100	100
82	o	80/87 (92%)	80 (100%)	0	100	100
83	p	135/181 (75%)	135 (100%)	0	100	100
84	q	142/178 (80%)	142 (100%)	0	100	100
85	r	147/169 (87%)	147 (100%)	0	100	100
86	s	340/381 (89%)	340 (100%)	0	100	100
87	t	40/158 (25%)	40 (100%)	0	100	100
87	u	31/158 (20%)	31 (100%)	0	100	100
87	v	31/158 (20%)	31 (100%)	0	100	100
87	w	30/158 (19%)	30 (100%)	0	100	100
87	x	30/158 (19%)	30 (100%)	0	100	100
87	y	30/158 (19%)	30 (100%)	0	100	100
88	H	182/228 (80%)	182 (100%)	0	100	100
89	z	226/287 (79%)	226 (100%)	0	100	100
All	All	13166/16790 (78%)	13166 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 115 such sidechains are listed below:

Mol	Chain	Res	Type
43	K	89	GLN
86	s	423	HIS
59	0	120	HIS
86	s	239	ASN
76	h	103	HIS

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	953/954 (99%)	136 (14%)	1 (0%)
32	Aw	67/68 (98%)	18 (26%)	0
33	Ax	68/70 (97%)	13 (19%)	0
34	Ay	68/70 (97%)	14 (20%)	0
35	Az	33/34 (97%)	14 (42%)	0
36	A	1556/1561 (99%)	224 (14%)	3 (0%)
37	B	71/72 (98%)	12 (16%)	0
All	All	2816/2829 (99%)	431 (15%)	4 (0%)

5 of 431 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	651	A
1	AA	673	U
1	AA	680	U
1	AA	688	A
1	AA	695	A

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	1539	C
36	A	2030	U
36	A	2112	A
36	A	2245	A

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
36	OMG	A	3040	36,32	23,26,27	0.22	0	32,38,41	0.36	0
53	AYA	U	2	53	6,7,8	0.83	0	6,8,10	0.55	0
29	AYA	A2	2	29	6,7,8	0.83	0	6,8,10	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	5F0	AI	184	9	8,8,9	0.55	0	8,9,11	1.10	1 (12%)
43	SAC	K	2	43	7,8,9	0.28	0	7,9,11	0.63	0
37	2MG	B	10	37	23,26,27	0.46	0	33,38,41	0.48	0
79	AYA	k	2	79	6,7,8	0.86	0	6,8,10	0.51	0
36	PSU	A	3067	36	18,21,22	0.77	0	21,30,33	2.78	6 (28%)
36	1MA	A	2617	36	21,25,26	0.75	1 (4%)	30,37,40	0.67	1 (3%)
1	5MC	AA	1488	1	19,22,23	0.28	0	26,32,35	0.41	0
37	PSU	B	39	37	18,21,22	0.76	0	21,30,33	2.76	5 (23%)
37	1MA	B	9	37	21,25,26	0.75	1 (4%)	30,37,40	0.66	1 (3%)
1	MA6	AA	1584	1	23,26,27	0.26	0	33,38,41	0.55	1 (3%)
36	OMU	A	3039	94,36	19,22,23	0.21	0	25,31,34	0.42	0
1	MA6	AA	1583	1	23,26,27	0.23	0	33,38,41	0.54	1 (3%)
17	AYA	AQ	2	17	6,7,8	0.82	0	6,8,10	0.50	0
36	OMG	A	2815	94,36,33	23,26,27	0.24	0	32,38,41	0.39	0
1	B8T	AA	1486	93,1	19,22,23	0.26	0	25,31,34	0.31	0
1	5MU	AA	1076	1	19,22,23	0.59	0	27,32,35	1.21	3 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
36	OMG	A	3040	36,32	-	0/9/27/28	0/3/3/3
53	AYA	U	2	53	-	0/5/6/8	-
29	AYA	A2	2	29	-	0/5/6/8	-
9	5F0	AI	184	9	-	0/9/9/10	-
43	SAC	K	2	43	-	0/7/8/10	-
37	2MG	B	10	37	-	0/9/27/28	0/3/3/3
79	AYA	k	2	79	-	1/5/6/8	-
36	PSU	A	3067	36	-	0/7/25/26	0/2/2/2
36	1MA	A	2617	36	-	0/7/25/26	0/3/3/3
1	5MC	AA	1488	1	-	0/7/25/26	0/2/2/2
37	PSU	B	39	37	-	0/7/25/26	0/2/2/2
37	1MA	B	9	37	-	0/7/25/26	0/3/3/3
1	MA6	AA	1584	1	-	2/11/29/30	0/3/3/3
36	OMU	A	3039	94,36	-	0/9/27/28	0/2/2/2
1	MA6	AA	1583	1	-	0/11/29/30	0/3/3/3
17	AYA	AQ	2	17	-	2/5/6/8	-
36	OMG	A	2815	94,36,33	-	0/9/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	B8T	AA	1486	93,1	-	0/7/27/28	0/2/2/2
1	5MU	AA	1076	1	-	0/7/25/26	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
37	B	9	1MA	C6-N6	3.08	1.35	1.28
36	A	2617	1MA	C6-N6	3.07	1.35	1.28

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
37	B	39	PSU	N1-C2-N3	8.62	124.26	115.17
36	A	3067	PSU	N1-C2-N3	8.61	124.25	115.17
37	B	39	PSU	C4-N3-C2	-6.86	116.92	126.37
36	A	3067	PSU	C4-N3-C2	-6.86	116.92	126.37
1	AA	1076	5MU	C4-N3-C2	-4.71	121.17	127.34

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
79	k	2	AYA	C-CA-N-CT
1	AA	1584	MA6	C5-C6-N6-C9
1	AA	1584	MA6	C4'-C5'-O5'-P
17	AQ	2	AYA	CB-CA-N-CT
17	AQ	2	AYA	C-CA-N-CT

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
36	A	3067	PSU	1	0
1	AA	1583	MA6	1	0
36	A	2815	OMG	1	0
1	AA	1486	B8T	2	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 278 ligands modelled in this entry, 265 are monoatomic - leaving 13 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
99	PUT	A	3304	-	5,5,5	0.23	0	4,4,4	0.50	0
92	SPD	A	3301	-	9,9,9	0.26	0	8,8,8	1.18	0
96	FES	AT	201	20,13	0,4,4	-	-	-	-	-
98	GDP	AX	503	-	29,30,30	0.42	0	45,47,47	0.43	0
92	SPD	A	3303	-	9,9,9	0.33	0	8,8,8	1.12	0
91	SPM	AA	1702	-	13,13,13	0.27	0	12,12,12	0.96	0
92	SPD	A	3302	-	9,9,9	0.32	0	8,8,8	1.17	0
97	ATP	AX	501	93	32,33,33	0.53	0	48,52,52	0.52	0
92	SPD	AA	1703	-	9,9,9	0.28	0	8,8,8	1.31	2 (25%)
100	VAL	B	101	37	4,6,7	0.54	0	6,7,9	0.96	0
96	FES	r	201	85,41	0,4,4	-	-	-	-	-
96	FES	AP	201	5,16	0,4,4	-	-	-	-	-
90	NAD	AA	1701	93	46,48,48	0.41	0	64,73,73	0.48	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
99	PUT	A	3304	-	-	0/3/3/3	-
92	SPD	A	3301	-	-	1/7/7/7	-
96	FES	AT	201	20,13	-	-	0/1/1/1
98	GDP	AX	503	-	-	0/16/32/32	0/3/3/3
92	SPD	A	3303	-	-	2/7/7/7	-
91	SPM	AA	1702	-	-	0/11/11/11	-
92	SPD	A	3302	-	-	0/7/7/7	-
97	ATP	AX	501	93	-	1/22/38/38	0/3/3/3
92	SPD	AA	1703	-	-	0/7/7/7	-
100	VAL	B	101	37	-	0/5/6/8	-
96	FES	r	201	85,41	-	-	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
96	FES	AP	201	5,16	-	-	0/1/1/1
90	NAD	AA	1701	93	-	0/30/62/62	0/5/5/5

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
92	AA	1703	SPD	C4-C5-N6	-2.23	106.09	112.07
92	AA	1703	SPD	C8-C7-N6	-2.08	106.49	112.07

There are no chirality outliers.

All (4) torsion outliers are listed below:

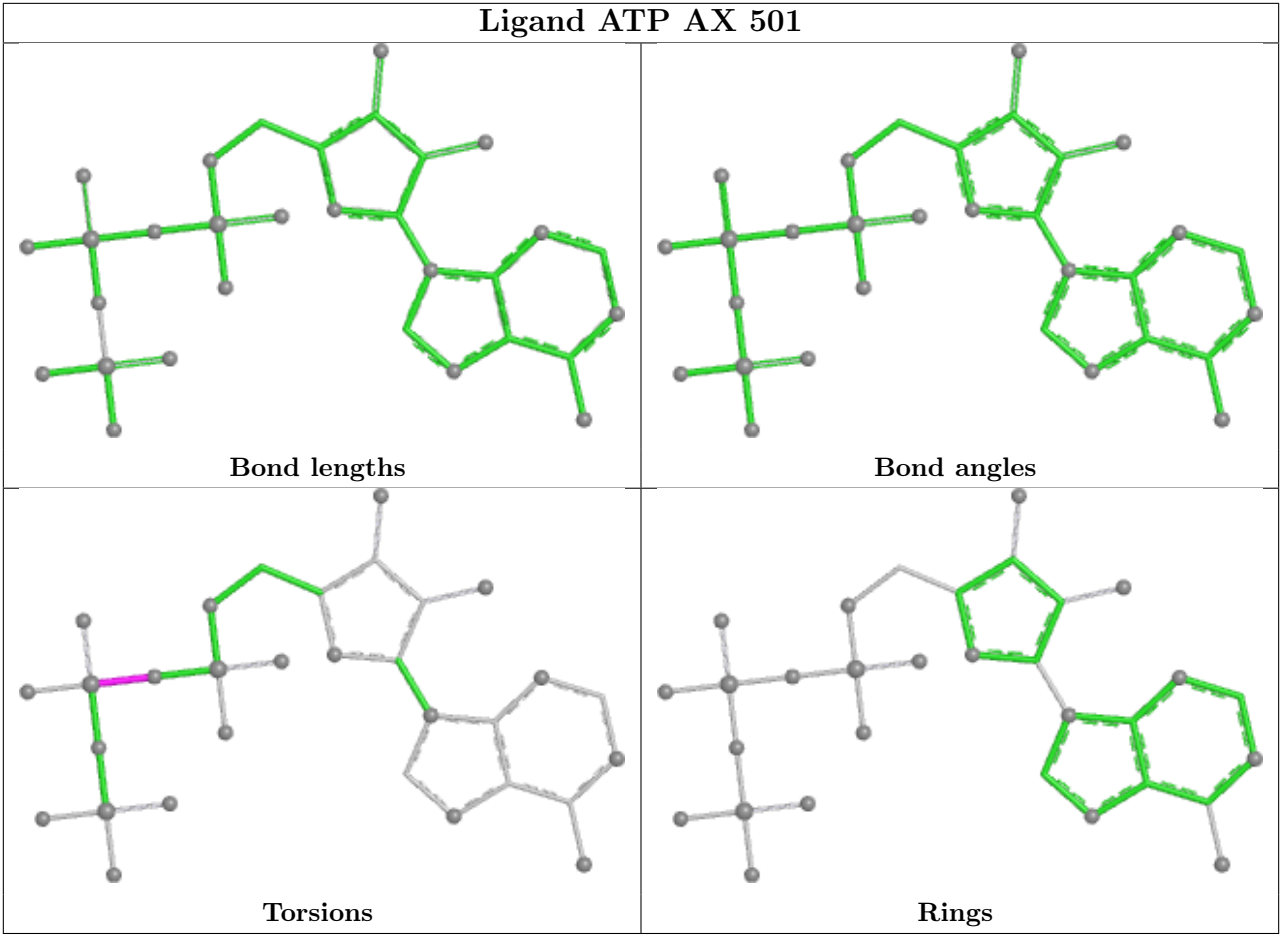
Mol	Chain	Res	Type	Atoms
92	A	3303	SPD	C4-C5-N6-C7
92	A	3301	SPD	N6-C7-C8-C9
92	A	3303	SPD	C8-C7-N6-C5
97	AX	501	ATP	PA-O3A-PB-O2B

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
90	AA	1701	NAD	1	0

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



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
34	Ay	1
33	Ax	1
35	Az	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	Ay	15:A	O3'	21:A	P	9.67
1	Ax	15:A	O3'	21:A	P	8.73
1	Az	3:A	O3'	4:A	P	3.00

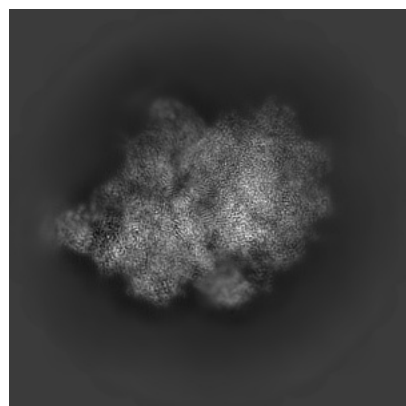
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13981. These allow visual inspection of the internal detail of the map and identification of artifacts.

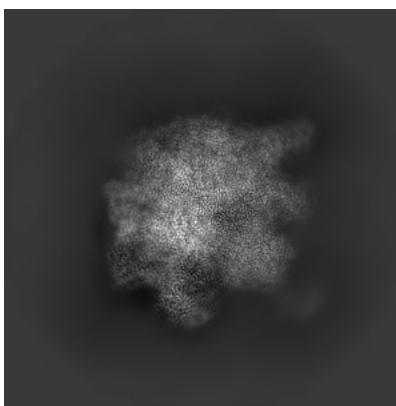
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

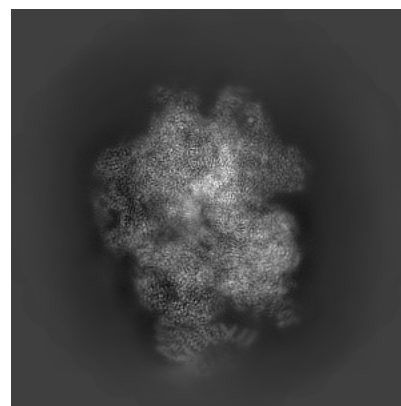
6.1.1 Primary map



X

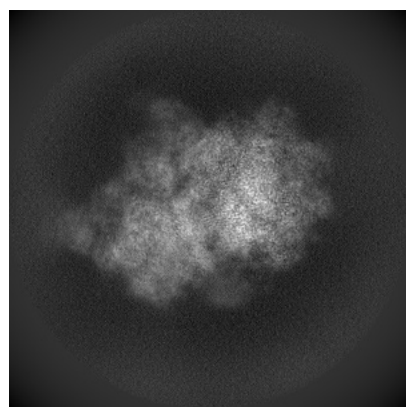


Y

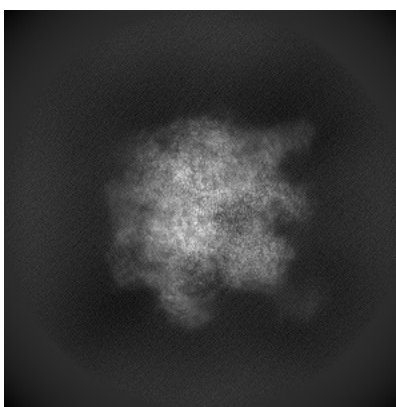


Z

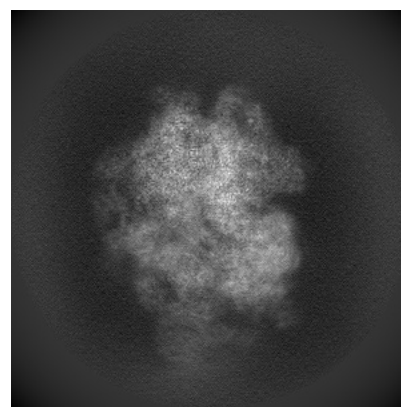
6.1.2 Raw map



X



Y

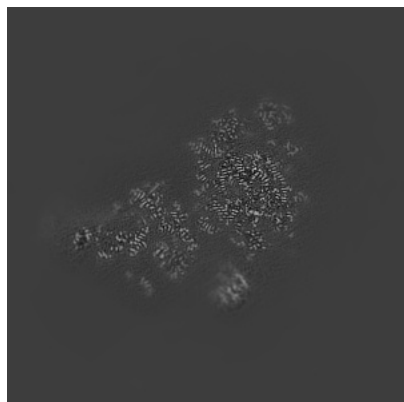


Z

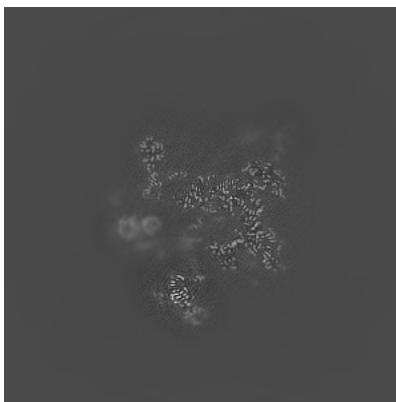
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

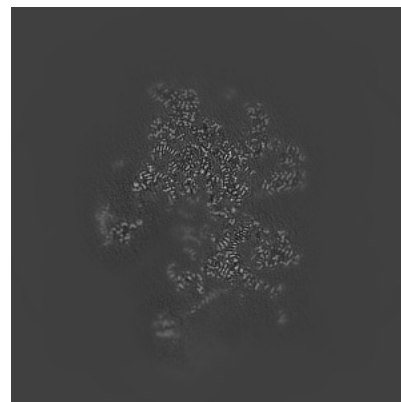
6.2.1 Primary map



X Index: 270

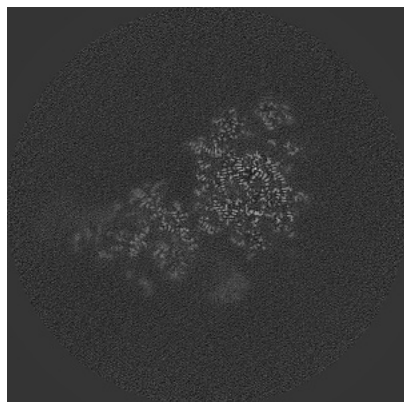


Y Index: 270

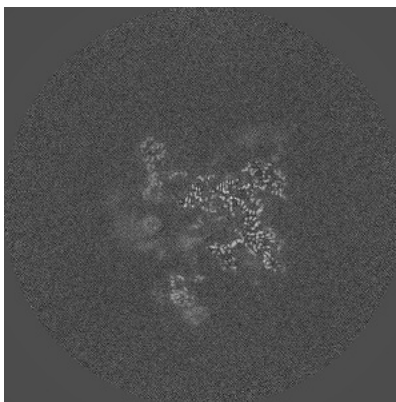


Z Index: 270

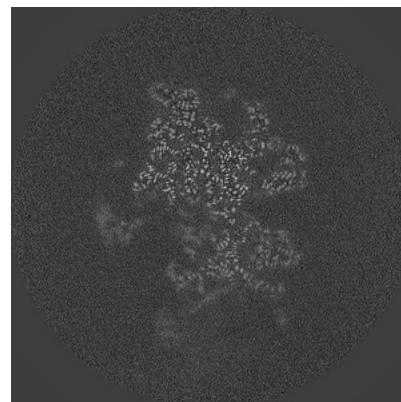
6.2.2 Raw map



X Index: 270



Y Index: 270

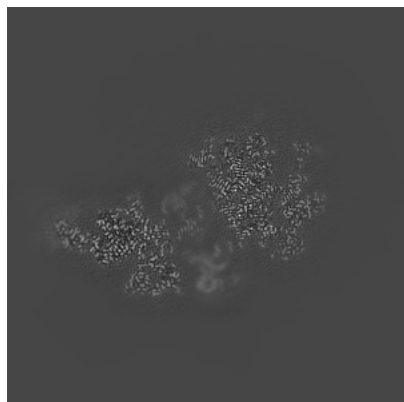


Z Index: 270

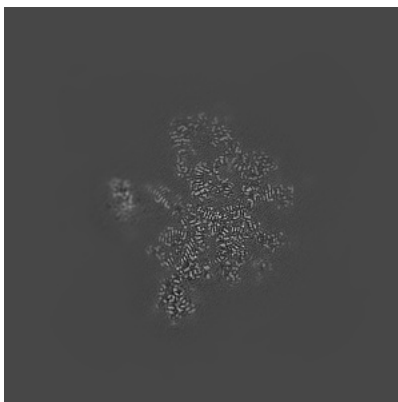
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

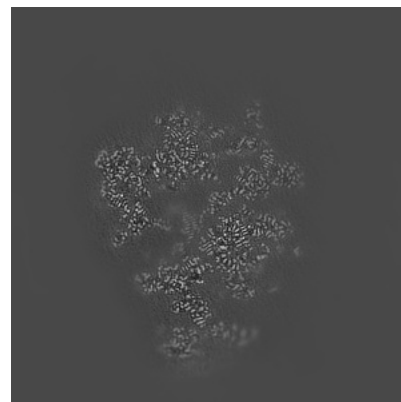
6.3.1 Primary map



X Index: 237

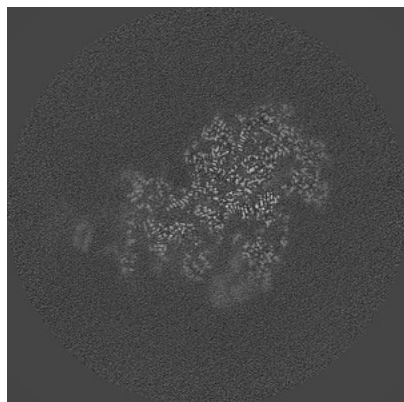


Y Index: 318

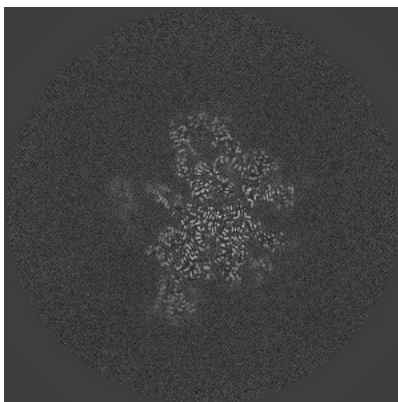


Z Index: 240

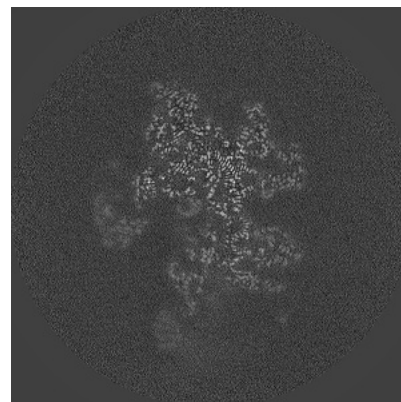
6.3.2 Raw map



X Index: 289



Y Index: 318

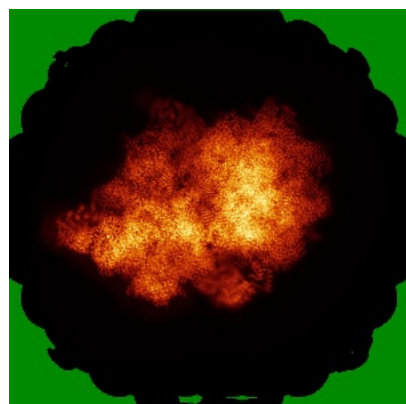


Z Index: 265

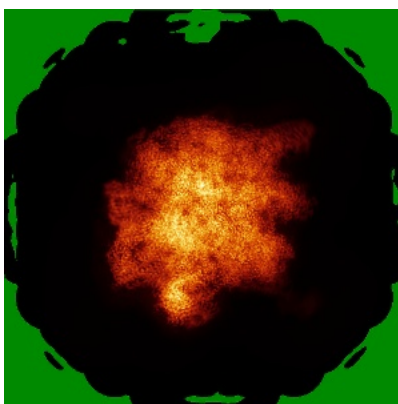
The images above show the largest variance slices of the map in three orthogonal directions.

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

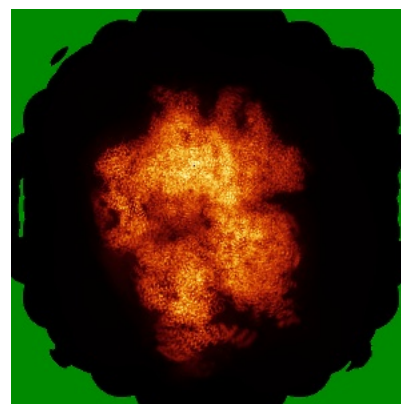
6.4.1 Primary map



X

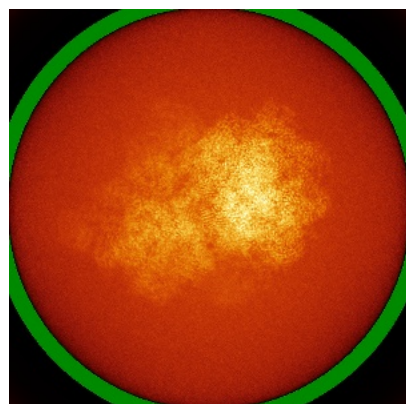


Y

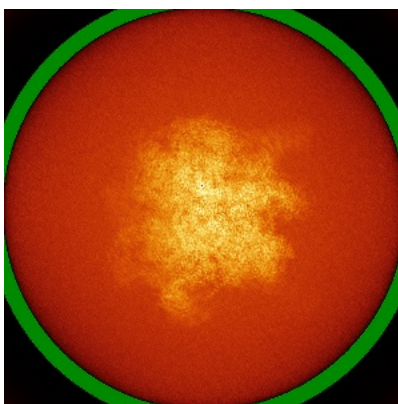


Z

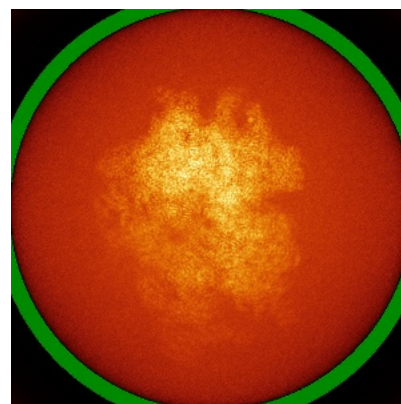
6.4.2 Raw map



X



Y



Z

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

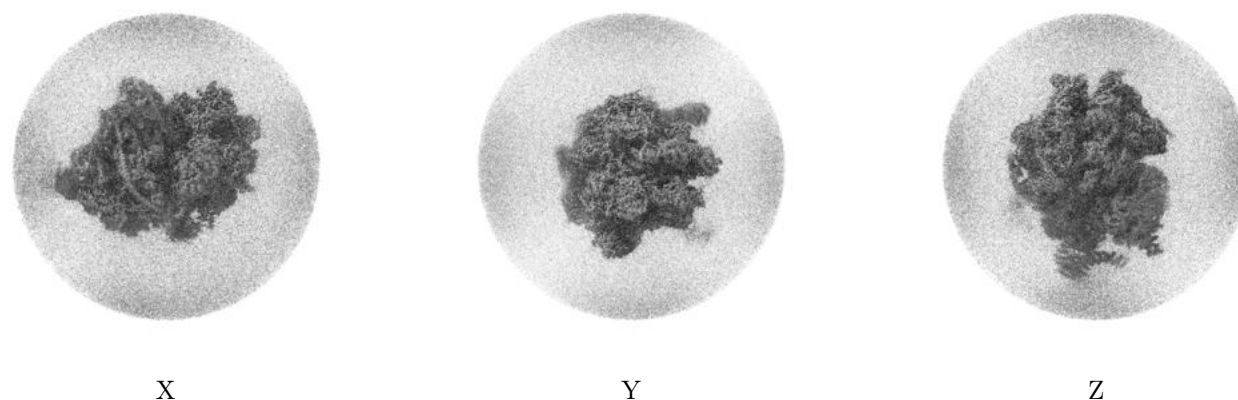
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 2.8. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

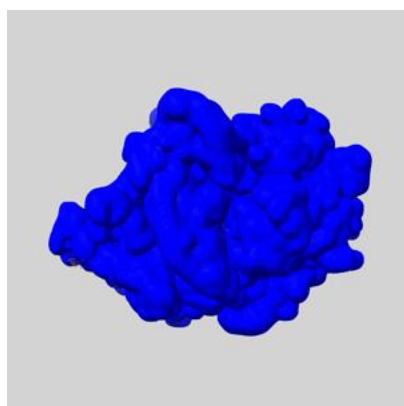
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

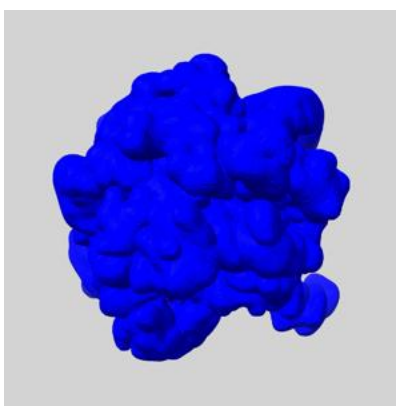
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

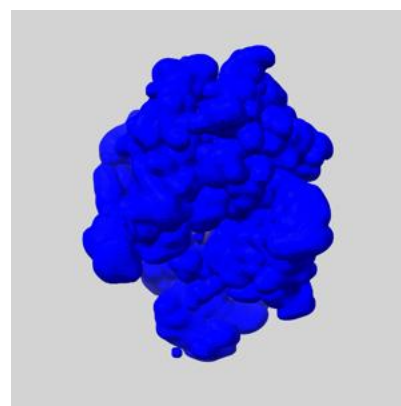
6.6.1 emd_13981_msk_1.map [i](#)



X

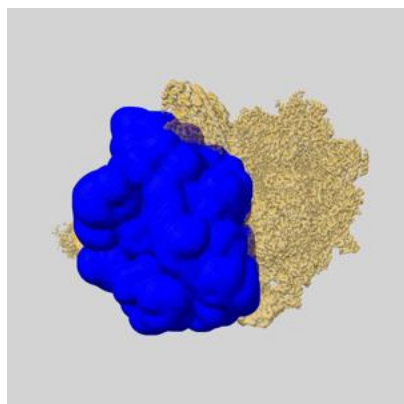


Y

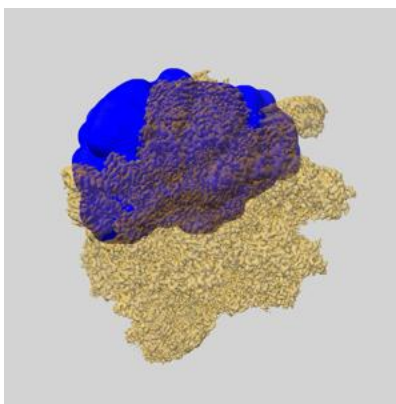


Z

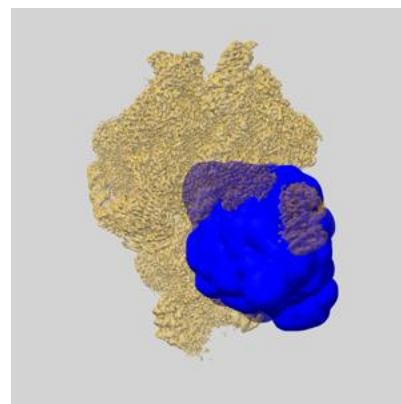
6.6.2 emd_13981_msk_2.map [i](#)



X

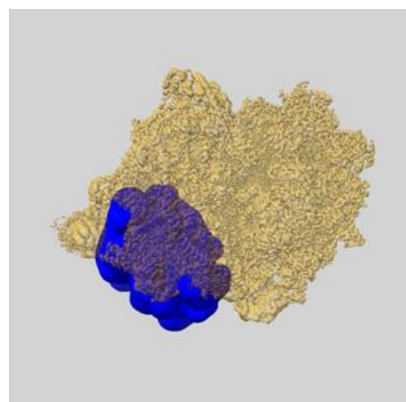


Y

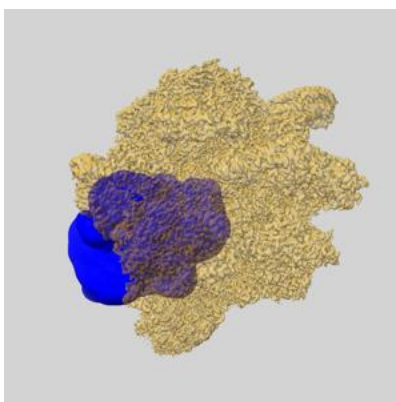


Z

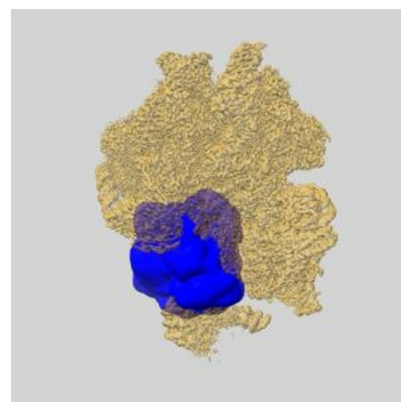
6.6.3 emd_13981_msk_3.map [i](#)



X

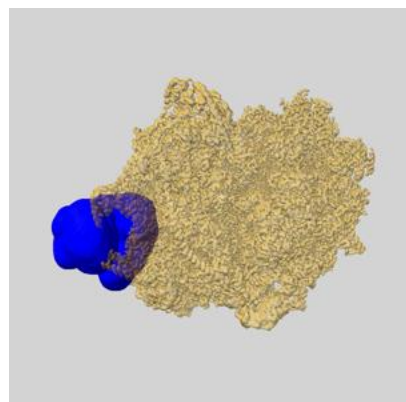


Y

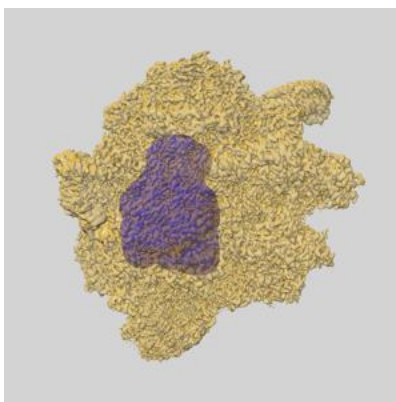


Z

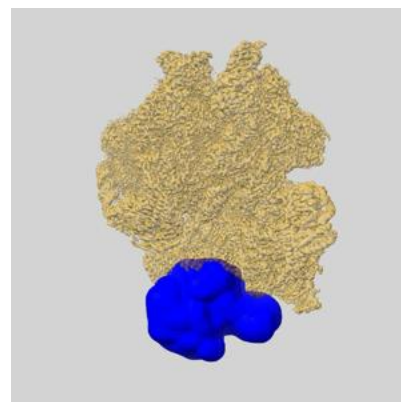
6.6.4 emd_13981_msk_4.map [i](#)



X

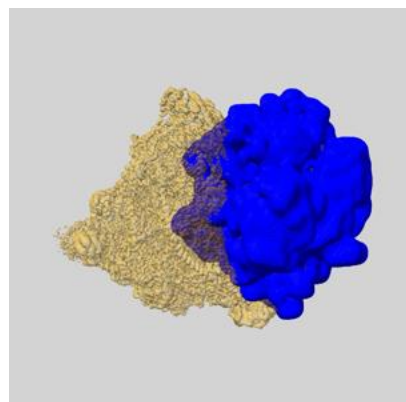


Y

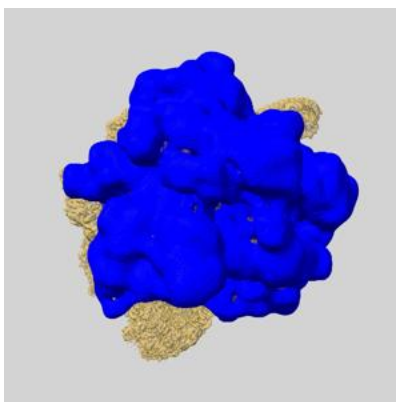


Z

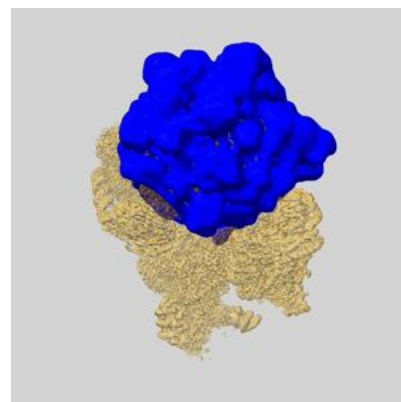
6.6.5 emd_13981_msk_5.map [i](#)



X

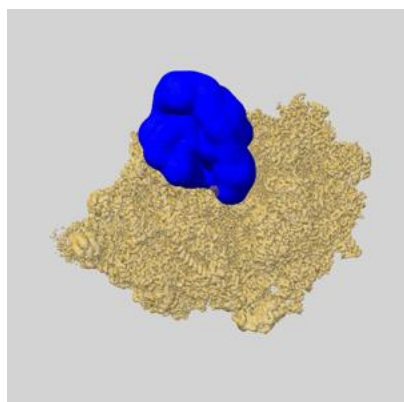


Y

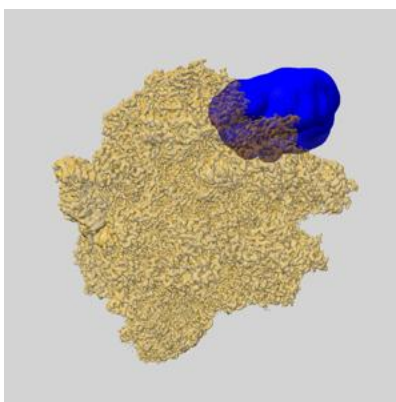


Z

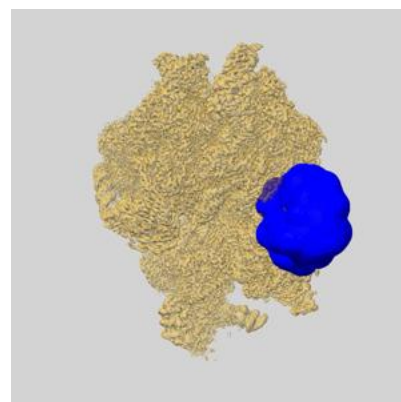
6.6.6 emd_13981_msk_6.map [i](#)



X

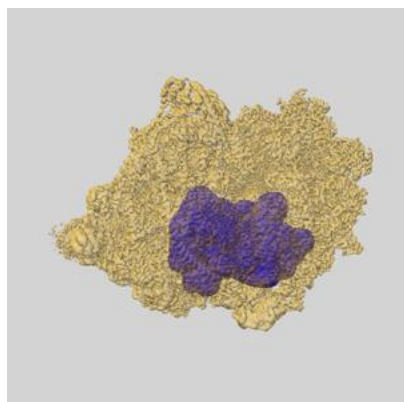


Y

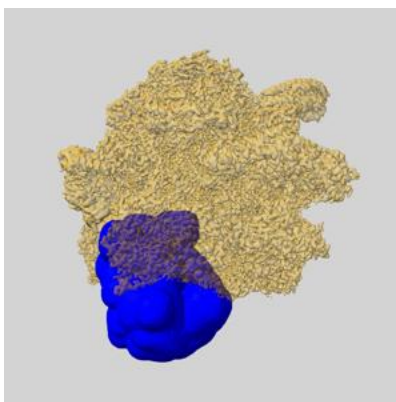


Z

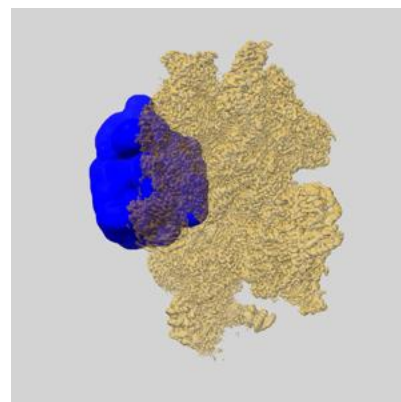
6.6.7 emd_13981_msk_7.map [i](#)



X

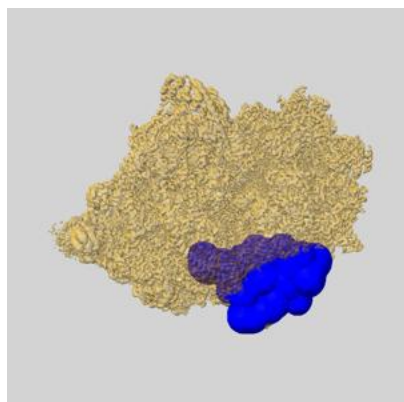


Y

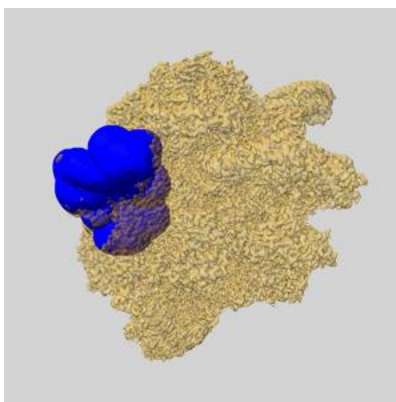


Z

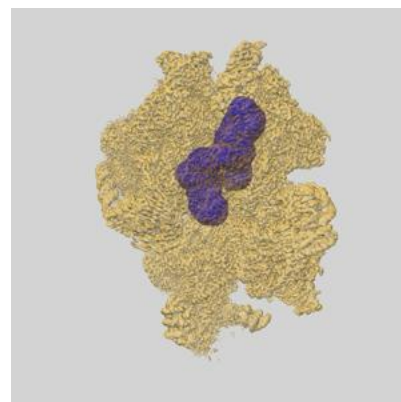
6.6.8 emd_13981_msk_8.map [i](#)



X

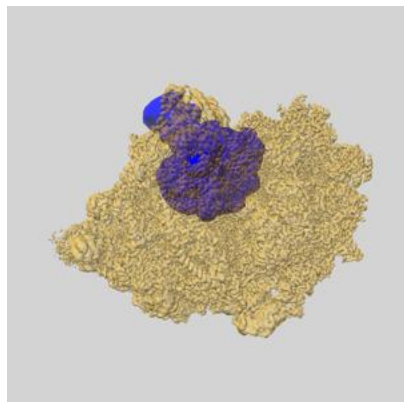


Y

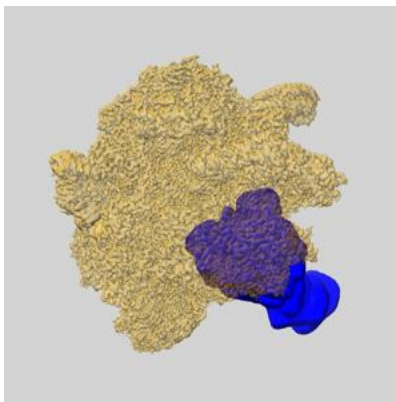


Z

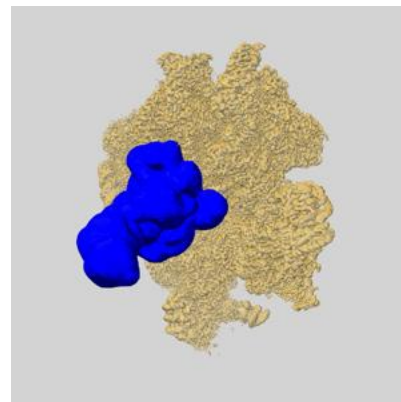
6.6.9 emd_13981_msk_9.map [i](#)



X



Y

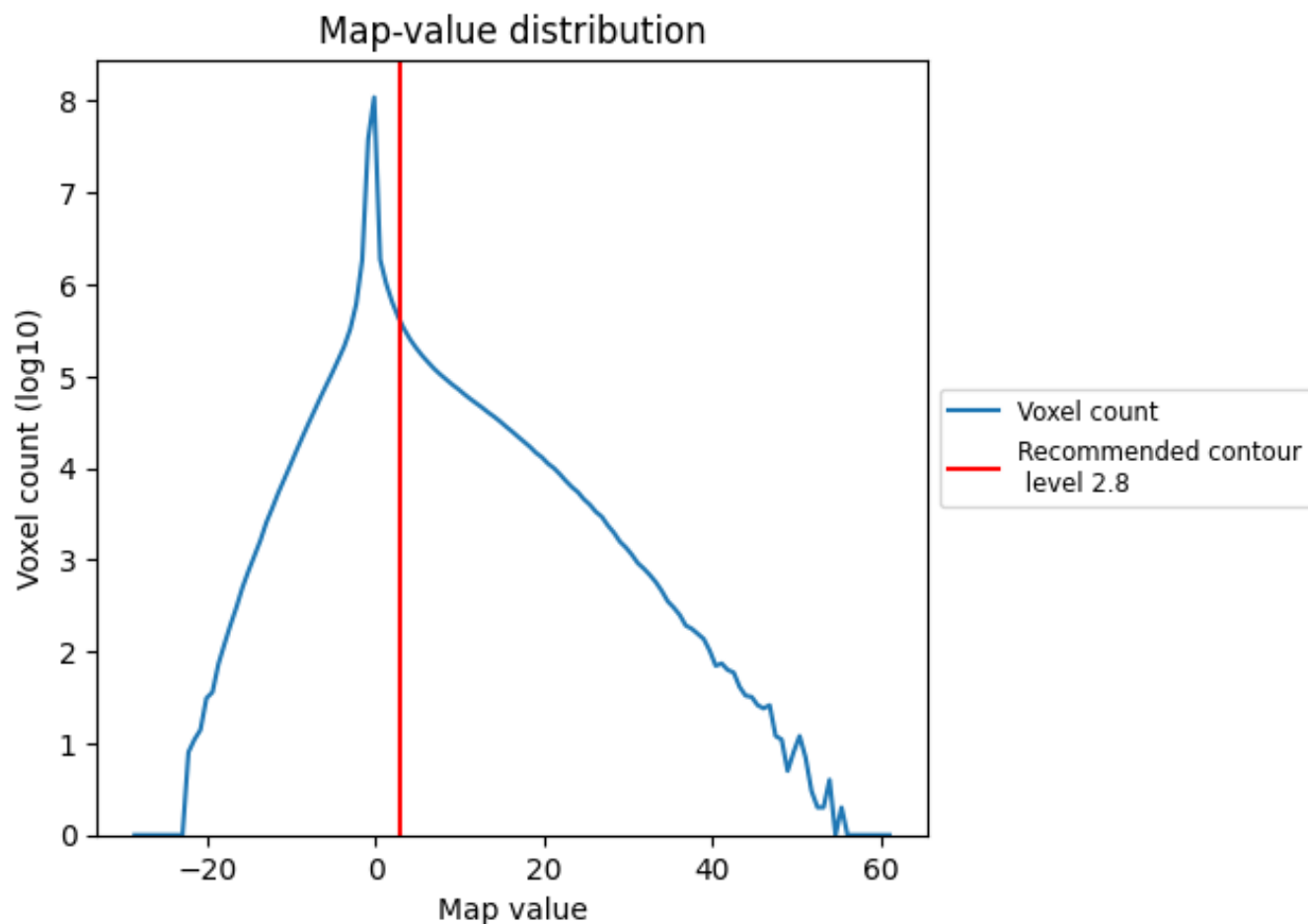


Z

7 Map analysis [i](#)

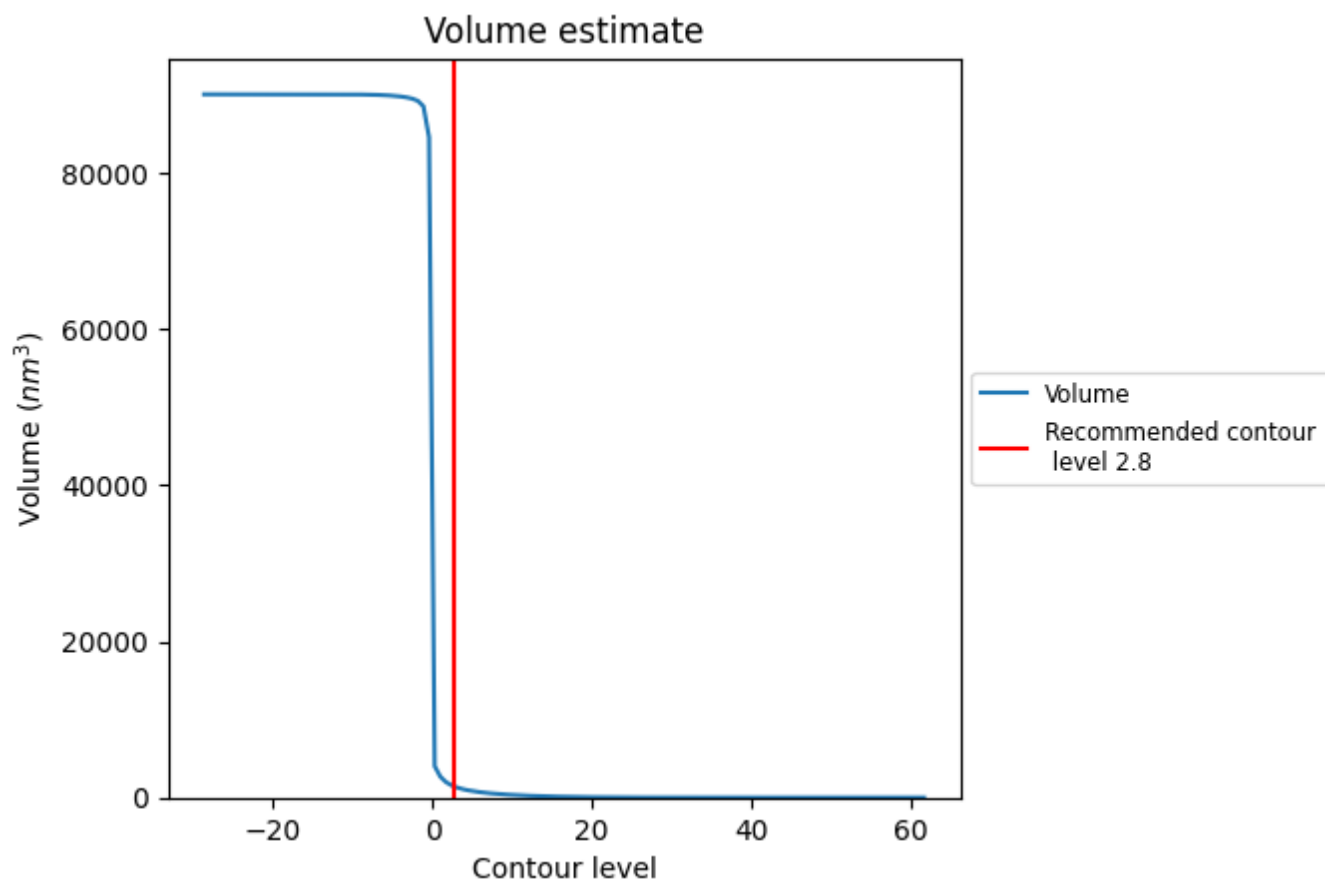
This section contains the results of statistical analysis of the map.

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.

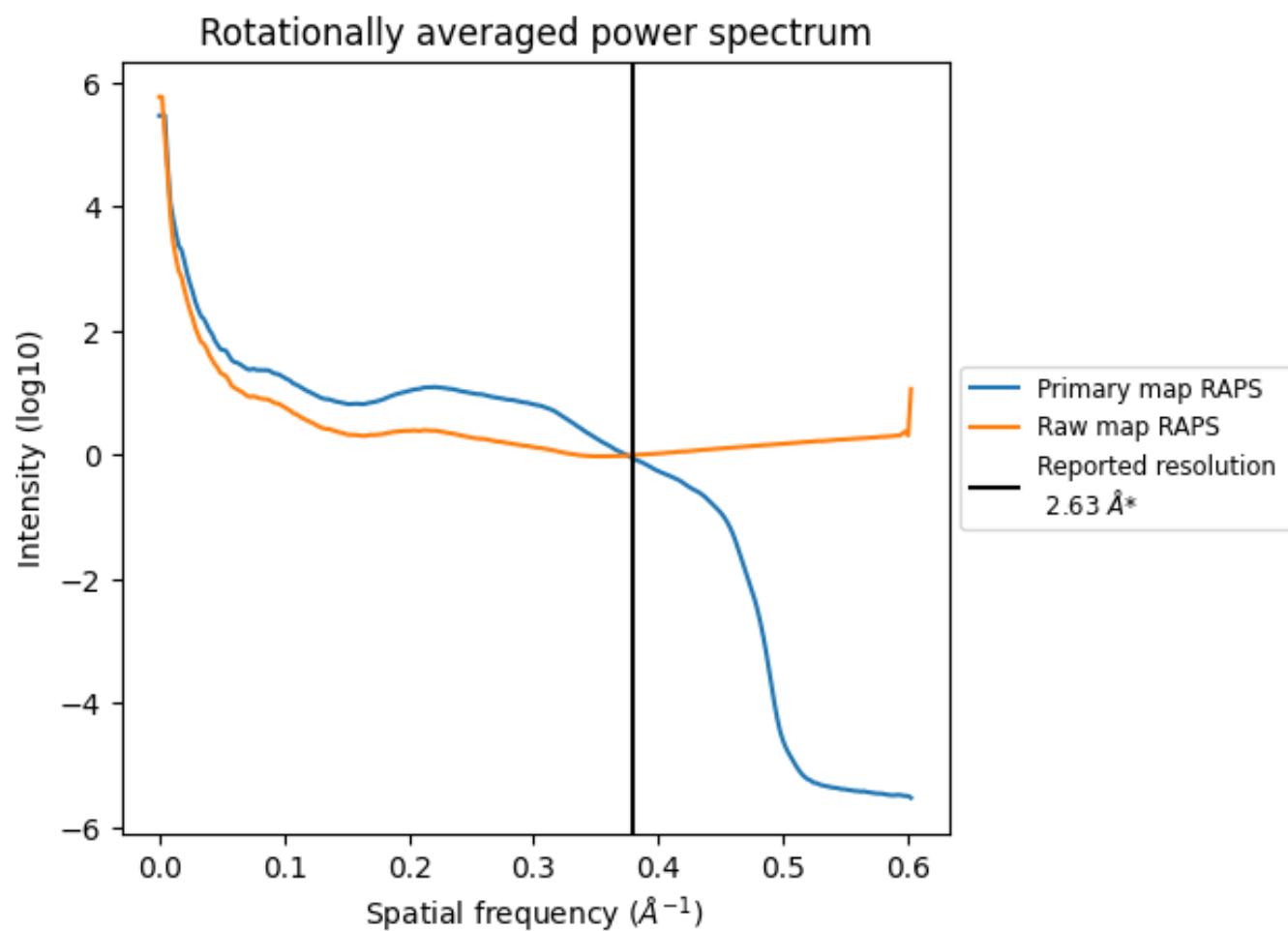
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1416 nm^3 ; this corresponds to an approximate mass of 1279 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 ⓘ

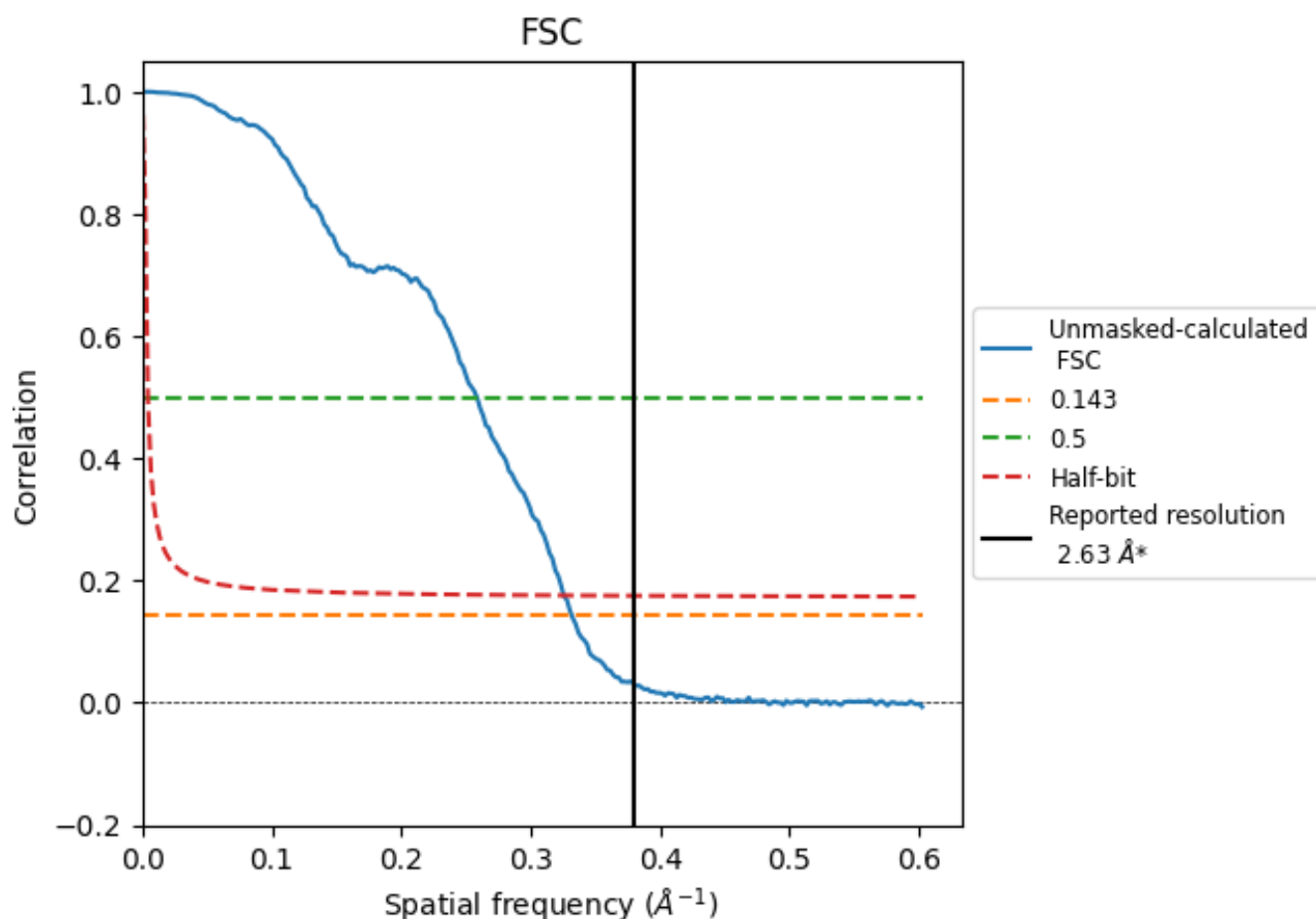


*Reported resolution corresponds to spatial frequency of 0.380 $\text{\AA}^{-1}$

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 0.380 Å⁻¹

8.2 Resolution estimates [i](#)

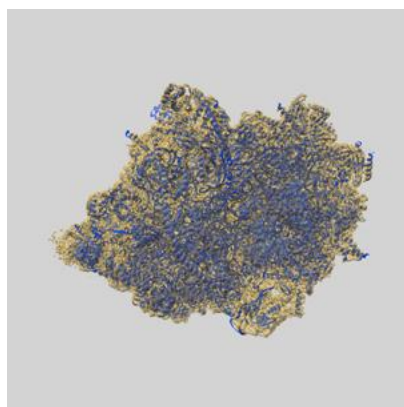
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.63	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.01	3.87	3.07

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.01 differs from the reported value 2.63 by more than 10 %

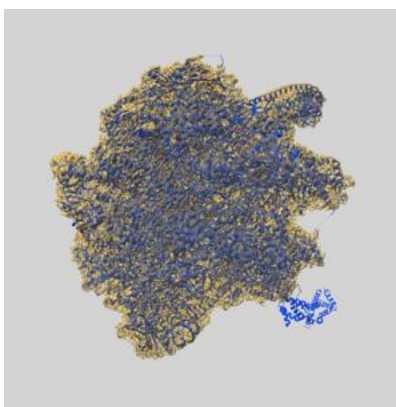
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13981 and PDB model 7QI5. Per-residue inclusion information can be found in section 3 on page 31.

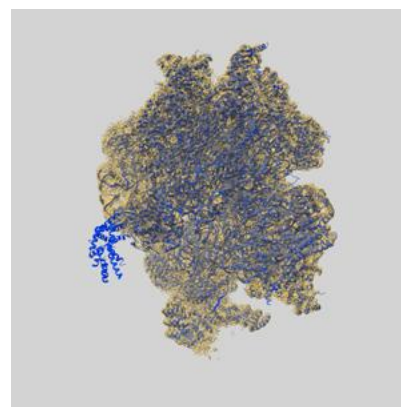
9.1 Map-model overlay [i](#)



X



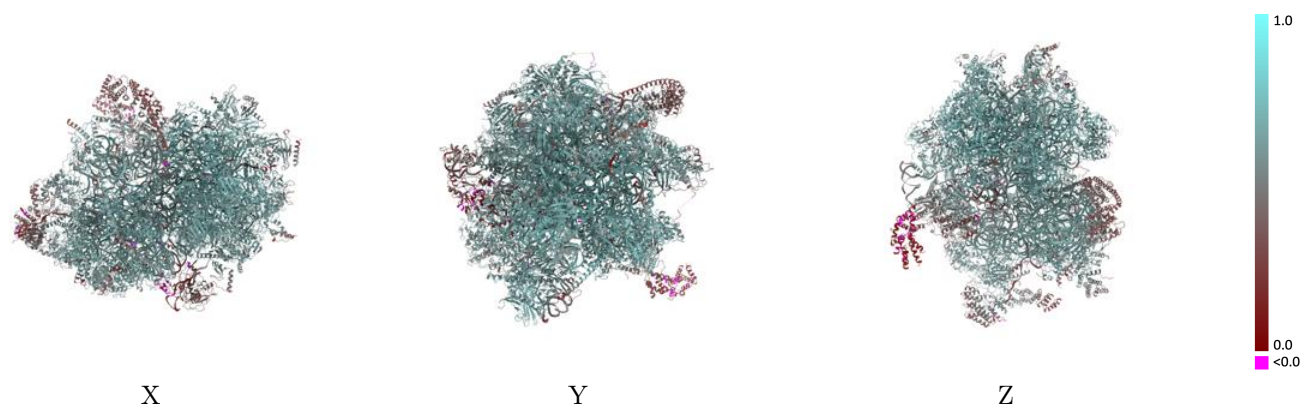
Y



Z

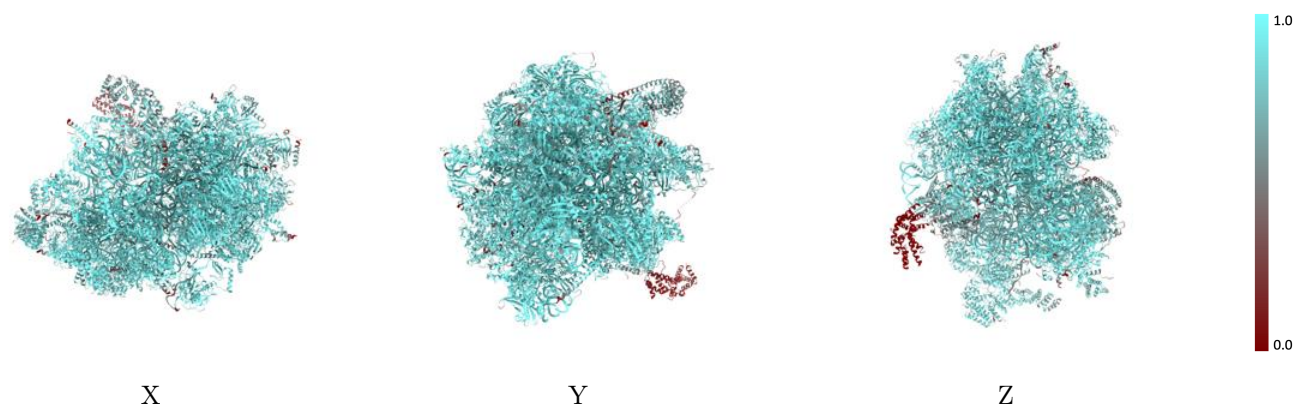
The images above show the 3D surface view of the map at the recommended contour level 2.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



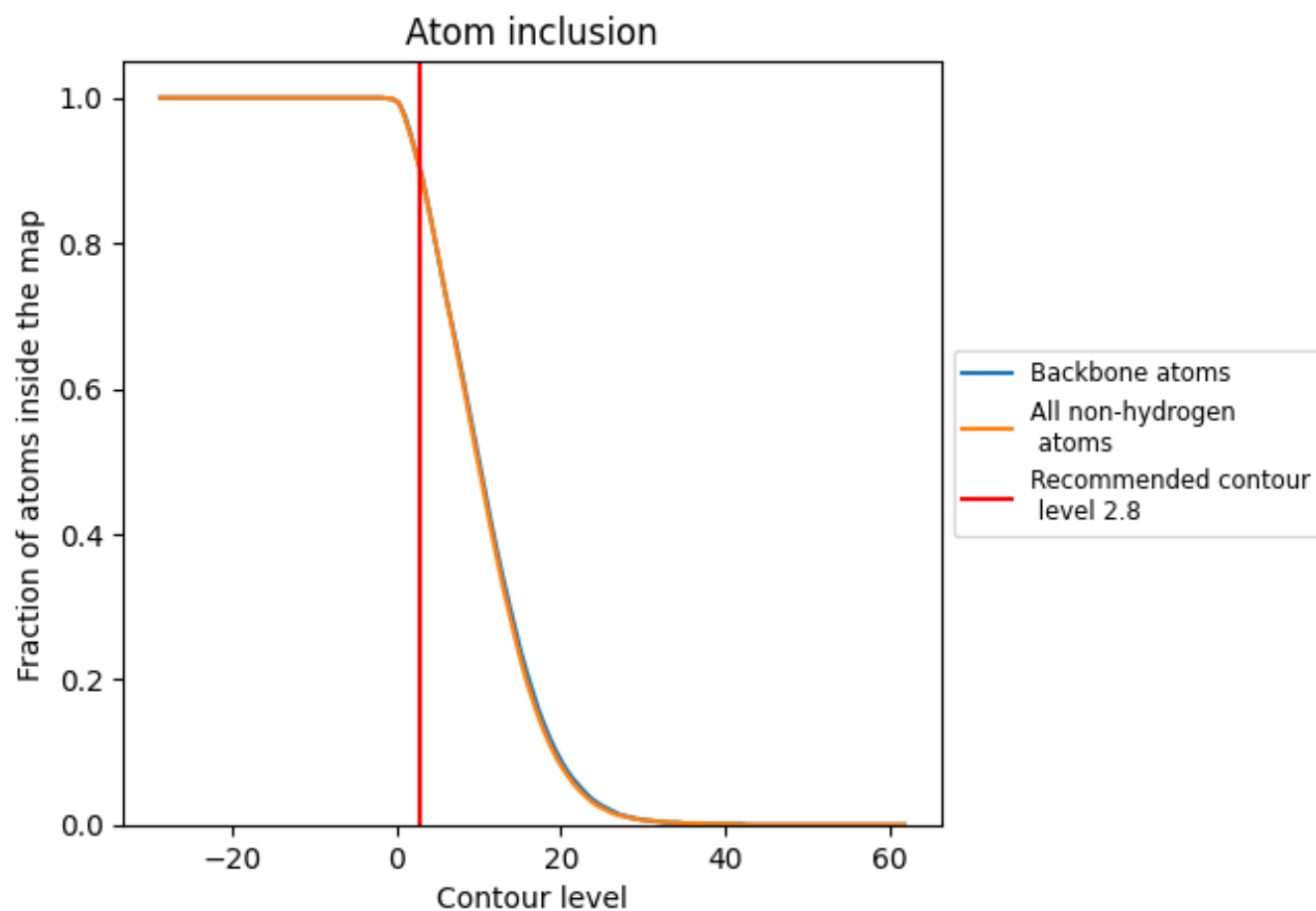
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.8).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 90% 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 (2.8) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9040	 0.5940
0	 0.9130	 0.6410
1	 0.8960	 0.6060
2	 0.9950	 0.7280
3	 0.9860	 0.7160
4	 0.9660	 0.6610
5	 0.9290	 0.6170
6	 0.9530	 0.6300
7	 0.8520	 0.5690
8	 0.8120	 0.5150
9	 0.8730	 0.5940
A	 0.9740	 0.6620
A0	 0.8550	 0.5140
A1	 0.9160	 0.5830
A2	 0.8650	 0.5540
A3	 0.9400	 0.6400
A4	 0.8730	 0.4560
AA	 0.9850	 0.6450
AB	 0.9400	 0.6180
AC	 0.9830	 0.6680
AD	 0.9130	 0.5920
AE	 0.9320	 0.6050
AF	 0.9390	 0.6000
AG	 0.8930	 0.5650
AH	 0.9540	 0.6210
AI	 0.9480	 0.6230
AJ	 0.9340	 0.6190
AK	 0.9820	 0.6700
AL	 0.8780	 0.5730
AM	 0.9220	 0.5940
AN	 0.9450	 0.6240
AO	 0.9220	 0.5860
AP	 0.9380	 0.6290
AQ	 0.9690	 0.6520
AR	 0.8710	 0.5340



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
AS	 0.8610	 0.5460
AT	 0.9190	 0.5990
AU	 0.8660	 0.5350
AV	 0.6510	 0.3600
AW	 0.9170	 0.5840
AX	 0.9500	 0.6050
AY	 0.8340	 0.5200
AZ	 0.9290	 0.6050
Aw	 0.7910	 0.4180
Ax	 0.8270	 0.4800
Ay	 0.9240	 0.3230
Az	 0.7050	 0.3920
B	 0.9690	 0.5600
D	 0.9640	 0.6670
E	 0.9470	 0.6590
F	 0.9670	 0.6800
H	 0.9110	 0.4640
I	 0.5630	 0.4350
J	 0.5080	 0.3640
K	 0.9680	 0.6860
L	 0.9550	 0.6610
M	 0.9590	 0.6740
N	 0.9310	 0.6530
O	 0.9460	 0.6600
P	 0.9790	 0.6660
Q	 0.8730	 0.6150
R	 0.9680	 0.6990
S	 0.9500	 0.6680
T	 0.9680	 0.6820
U	 0.8550	 0.6020
V	 0.8480	 0.5720
W	 0.9770	 0.6850
X	 0.9070	 0.6100
Y	 0.9380	 0.6460
Z	 0.9430	 0.6700
a	 0.8450	 0.5740
b	 0.9620	 0.6690
c	 0.9000	 0.6090
d	 0.7810	 0.5240
e	 0.8530	 0.5230
f	 0.8710	 0.5550
g	 0.9290	 0.6400

Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
h	 0.8450	 0.5610
i	 0.9770	 0.7080
j	 0.8710	 0.5990
k	 0.7530	 0.5080
l	 0.5990	 0.4290
m	 0.7940	 0.4870
o	 0.9700	 0.6960
p	 0.8070	 0.5350
q	 0.6890	 0.4570
r	 0.9260	 0.6280
s	 0.9350	 0.6330
t	 0.0680	 0.2330
u	 0.0000	 0.2200
v	 0.0000	 0.1390
w	 0.0000	 0.0470
x	 0.0000	 0.0870
y	 0.0000	 0.0550
z	 0.8680	 0.2230