



wwPDB EM Validation Summary Report ⓘ

Nov 20, 2022 – 12:20 pm GMT

PDB ID : 6I9R
EMDB ID : EMD-4434
Title : Large subunit of the human mitochondrial ribosome in complex with Virginiamycin M and Quinupristin
Authors : Modelska, A.; Aibara, S.; Amunts, A.
Deposited on : 2018-11-25
Resolution : 3.90 Å(reported)

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

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

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

The types of validation reports are described at
<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

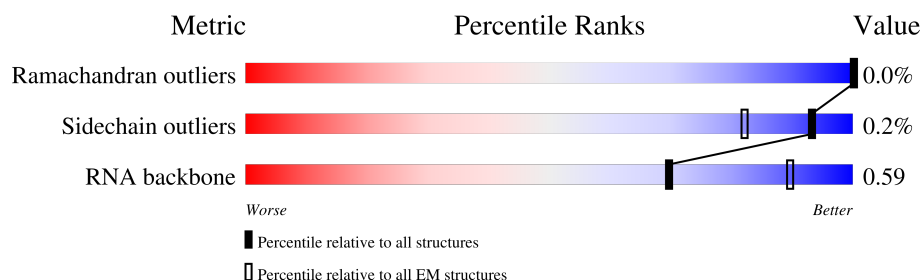
EMDB validation analysis : 0.0.1.dev43
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

1 Overall quality at a glance

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

The reported resolution of this entry is 3.90 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	0	188	
2	1	65	
3	2	92	
4	3	188	
5	4	103	
6	5	423	
7	6	380	
8	7	338	



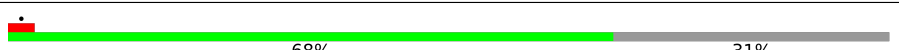
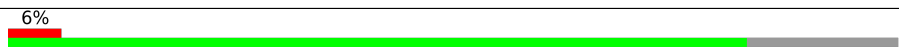

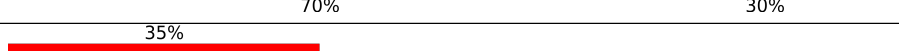
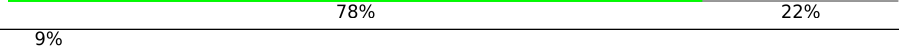





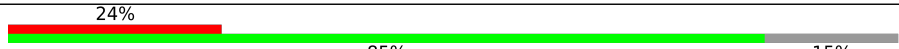


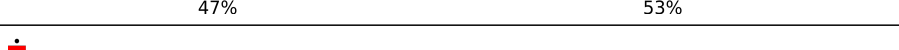
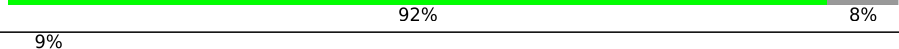


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Mol	Chain	Length	Quality of chain
9	8	206	
10	9	137	
11	A	1557	
12	B	69	
13	D	305	
14	E	348	
15	F	311	
16	H	267	
17	I	261	
18	J	192	
19	K	178	
20	L	145	
21	M	296	
22	N	251	
23	O	175	
24	P	179	
25	Q	292	
26	R	149	
27	S	205	
28	T	212	
29	U	153	
30	V	216	
31	W	148	
32	X	256	
33	Y	250	

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Mol	Chain	Length	Quality of chain
34	Z	161	
35	a	142	
36	b	215	
37	c	332	
38	d	306	
39	e	279	
40	f	212	
41	g	166	
42	h	158	
43	i	128	
44	j	123	
45	k	112	
46	l	138	
47	m	128	
48	o	102	
49	p	206	
50	q	222	
51	r	196	
52	s	439	

2 Entry composition

There are 57 unique types of molecules in this entry. The entry contains 100274 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 39S ribosomal protein L32, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	108	Total	C	N	O	S	0	0
			880	545	172	157	6		

- Molecule 2 is a protein called 39S ribosomal protein L33, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	53	Total	C	N	O	S	0	0
			439	281	84	72	2		

- Molecule 3 is a protein called 39S ribosomal protein L34, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	46	Total	C	N	O	S	0	0
			376	233	83	59	1		

- Molecule 4 is a protein called 39S ribosomal protein L35, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	3	95	Total	C	N	O	S	0	0
			831	539	162	127	3		

- Molecule 5 is a protein called 39S ribosomal protein L36, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	4	38	Total	C	N	O	S	0	0
			341	217	72	48	4		

- Molecule 6 is a protein called 39S ribosomal protein L37, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	5	393	Total	C	N	O	S	0	0
			3204	2070	559	564	11		

- Molecule 7 is a protein called 39S ribosomal protein L38, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	6	354	Total	C	N	O	S	0	0
			2947	1881	525	532	9		

- Molecule 8 is a protein called 39S ribosomal protein L39, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	7	291	Total	C	N	O	S	0	0
			2365	1514	401	432	18		

- Molecule 9 is a protein called 39S ribosomal protein L40, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	8	105	Total	C	N	O	S	0	0
			886	568	152	164	2		

- Molecule 10 is a protein called 39S ribosomal protein L41, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	9	124	Total	C	N	O	S	0	0
			996	644	170	180	2		

- Molecule 11 is a RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	A	1501	Total	C	N	O	P	0	0
			31875	14306	5758	10310	1501		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2360	A	U	conflict	GB 1391841121
A	2361	C	G	conflict	GB 1391841121

- Molecule 12 is a RNA chain called tRNA Val.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	B	60	Total	C	N	O	P	0	0
			1275	572	230	413	60		

- Molecule 13 is a protein called 39S ribosomal protein L2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	D	236	Total	C	N	O	S	0	0
			1842	1145	373	315	9		

- Molecule 14 is a protein called 39S ribosomal protein L3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	E	304	Total	C	N	O	S	0	0
			2396	1539	416	430	11		

- Molecule 15 is a protein called 39S ribosomal protein L4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	F	250	Total	C	N	O	S	0	0
			2013	1294	365	348	6		

- Molecule 16 is a protein called 39S ribosomal protein L9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	H	95	Total	C	N	O		0	0
			784	498	152	134			

- Molecule 17 is a protein called 39S ribosomal protein L10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	I	167	Total	C	N	O	S	0	0
			1350	869	245	226	10		

- Molecule 18 is a protein called 39S ribosomal protein L11, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	J	170	Total	C	N	O	S	0	0
			1294	826	230	236	2		

- Molecule 19 is a protein called 39S ribosomal protein L13, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	K	177	Total	C	N	O	S	0	0
			1451	934	259	251	7		

- Molecule 20 is a protein called 39S ribosomal protein L14, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	L	115	Total	C	N	O	S	0	0
			889	559	171	154	5		

- Molecule 21 is a protein called 39S ribosomal protein L15, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	M	287	Total	C	N	O	S	0	0
			2305	1472	425	402	6		

- Molecule 22 is a protein called 39S ribosomal protein L16, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	N	221	Total	C	N	O	S	0	0
			1778	1138	325	305	10		

- Molecule 23 is a protein called 39S ribosomal protein L17, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	O	152	Total	C	N	O	S	0	0
			1245	784	239	215	7		

- Molecule 24 is a protein called Mitochondrial ribosomal protein L18, isoform CRA_b.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	P	143	Total	C	N	O	S	0	0
			1164	729	223	207	5		

- Molecule 25 is a protein called 39S ribosomal protein L19, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	Q	219	Total	C	N	O	S	0	0
			1822	1168	322	323	9		

- Molecule 26 is a protein called 39S ribosomal protein L20, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	R	140	Total	C	N	O	S	0	0
			1153	732	231	186	4		

- Molecule 27 is a protein called 39S ribosomal protein L21, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	S	160	Total	C	N	O	S	0	0
			1284	829	226	225	4		

- Molecule 28 is a protein called 39S ribosomal protein L22, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	T	166	Total	C	N	O	S	0	0
			1368	875	254	232	7		

- Molecule 29 is a protein called 39S ribosomal protein L23, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	U	141	Total	C	N	O	S	0	0
			1171	743	222	203	3		

- Molecule 30 is a protein called 39S ribosomal protein L24, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	V	202	Total	C	N	O	S	0	0
			1652	1053	294	297	8		

- Molecule 31 is a protein called 39S ribosomal protein L27, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	W	111	Total	C	N	O	S	0	0
			871	558	164	146	3		

- Molecule 32 is a protein called 39S ribosomal protein L28, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	X	243	Total	C	N	O	S	0	0
			2027	1310	350	362	5		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	148	ALA	THR	conflict	UNP Q13084
X	149	SER	PRO	conflict	UNP Q13084
X	150	GLY	LYS	conflict	UNP Q13084

- Molecule 33 is a protein called 39S ribosomal protein L47, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	Y	178	Total	C	N	O	S	0	0
			1534	981	295	254	4		

- Molecule 34 is a protein called 39S ribosomal protein L30, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	Z	120	Total	C	N	O	S	0	0
			978	626	183	166	3		

- Molecule 35 is a protein called 39S ribosomal protein L42, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	a	83	Total	C	N	O	S	0	0
			693	440	124	124	5		

- Molecule 36 is a protein called 39S ribosomal protein L43, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	b	148	Total	C	N	O	S	0	0
			1178	733	229	213	3		

- Molecule 37 is a protein called 39S ribosomal protein L44, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	c	275	Total	C	N	O	S	0	0
			2217	1415	383	410	9		

- Molecule 38 is a protein called 39S ribosomal protein L45, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	d	214	Total	C	N	O	S	0	0
			1736	1113	303	307	13		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
d	62	PHE	GLU	conflict	UNP Q9BRJ2
d	63	ALA	PHE	conflict	UNP Q9BRJ2

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

Mol	Chain	Residues	Atoms					AltConf	Trace
39	e	217	Total	C	N	O	S	0	0
			1762	1124	310	323	5		

- Molecule 40 is a protein called 39S ribosomal protein L48, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	f	125	Total	C	N	O	S	0	0
			999	640	159	196	4		

- Molecule 41 is a protein called 39S ribosomal protein L49, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	g	131	Total	C	N	O	S	0	0
			1085	701	190	192	2		

- Molecule 42 is a protein called 39S ribosomal protein L50, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	h	109	Total	C	N	O	S	0	0
			886	562	155	166	3		

- Molecule 43 is a protein called 39S ribosomal protein L51, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	i	97	Total	C	N	O	S	0	0
			827	532	165	126	4		

- Molecule 44 is a protein called cDNA FLJ76418, highly similar to Homo sapiens mitochondrial ribosomal protein L52 (MRPL52), transcript variant 1, mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	j	86	Total	C	N	O	S	0	0
			689	426	134	127	2		

- Molecule 45 is a protein called 39S ribosomal protein L53, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	k	95	Total	C	N	O	S	0	0
			732	456	139	132	5		

- Molecule 46 is a protein called 39S ribosomal protein L54, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	l	80	Total	C	N	O	S	0	0
			675	428	118	126	3		

- Molecule 47 is a protein called 39S ribosomal protein L55, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	m	60	Total	C	N	O	S	0	0
			499	309	103	85	2		

- Molecule 48 is a protein called Ribosomal protein 63, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	o	94	Total	C	N	O	S	0	0
			797	501	165	128	3		

- Molecule 49 is a protein called Peptidyl-tRNA hydrolase ICT1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	p	127	Total	C	N	O	S	0	0
			1058	661	201	192	4		

- Molecule 50 is a protein called Growth arrest and DNA damage-inducible proteins-interacting protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	q	130	Total	C	N	O	S	0	0
			1092	680	213	194	5		

- Molecule 51 is a protein called 39S ribosomal protein S18a, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	r	153	Total	C	N	O	S	0	0
			1256	797	241	210	8		

- Molecule 52 is a protein called 39S ribosomal protein S30, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	s	370	Total	C	N	O	S	0	0
			3036	1946	542	534	14		

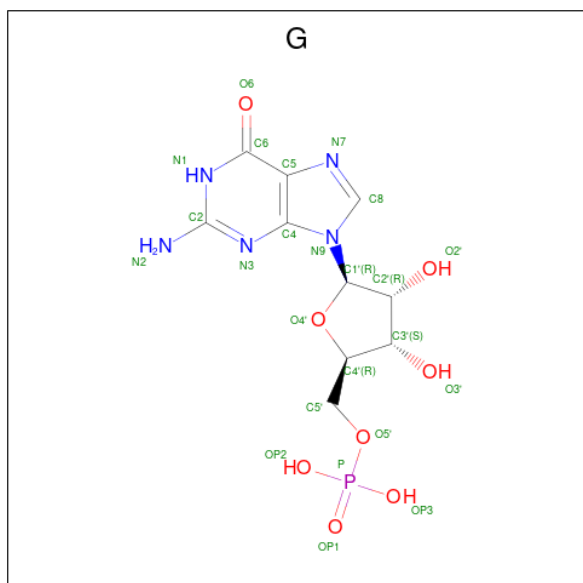
- Molecule 53 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
53	0	1	Total	Zn	0
			1	1	
53	4	1	Total	Zn	0
			1	1	
53	I	1	Total	Zn	0
			1	1	

- Molecule 54 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

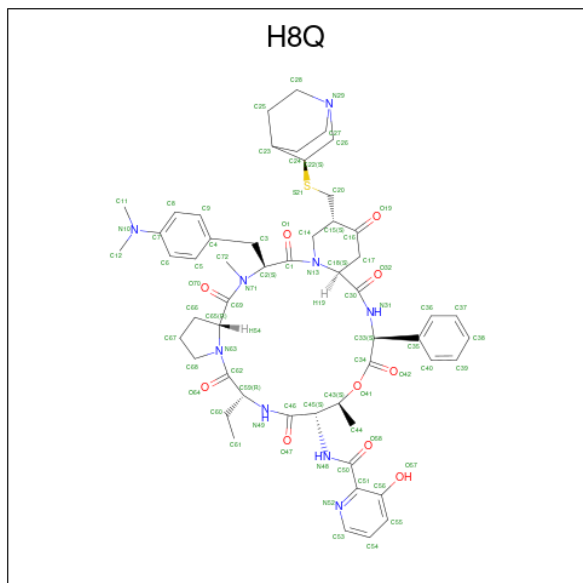
Mol	Chain	Residues	Atoms		AltConf
54	9	1	Total	Mg	0
			1	1	
54	A	128	Total	Mg	0
			128	128	
54	D	1	Total	Mg	0
			1	1	
54	L	1	Total	Mg	0
			1	1	
54	M	1	Total	Mg	0
			1	1	
54	g	1	Total	Mg	0
			1	1	
54	o	1	Total	Mg	0
			1	1	

- Molecule 55 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: G) (formula: $C_{10}H_{14}N_5O_8P$).



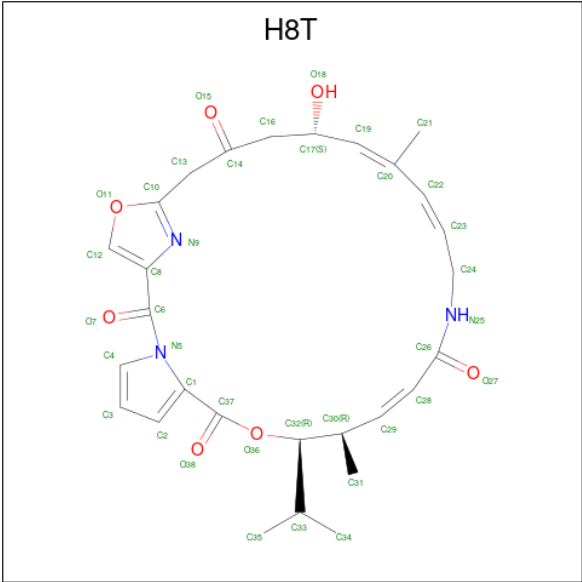
Mol	Chain	Residues	Atoms					AltConf
55	A	1	Total	C	N	O	P	0
			23	10	5	7	1	

- Molecule 56 is {N}-[(3 {S},6 {R},12 {R},15 {S},16 {S},19 {S},22 {S},25 {S})-25-[[[(3 {S})-1-azabicyclo[2.2.2]octan-3-yl]sulfanylmethyl]-3-[[4-(dimethylamino)phenyl]methyl]-12-ethyl-4,16-dimethyl-2,5,11,14,18,21,24-heptakis(oxidanylidene)-19-phenyl-17-oxa-1,4,10,13,20-pentazatricyclo[20.4.0.0[^]{6,10}]hexacosan-15-yl]-3-oxidanyl-pyridine-2-carboxamide (three-letter code: H8Q) (formula: C₅₃H₆₇N₉O₁₀S).



Mol	Chain	Residues	Atoms					AltConf
56	A	1	Total	C	N	O	S	0
			73	53	9	10	1	

- Molecule 57 is (10 {R},11 {R},12 {E},17 {E},19 {E},21 {S})-11,19-dimethyl-21-oxidanyl-10-propan-2-yl-9,26-dioxa-3,15,28-triazatricyclo[23.2.1.0[^]{3,7}]octacosan-1(27),4,6,12,17,19,25(28)-heptaene-2,8,14,23-tetrone (three-letter code: H8T) (formula: C₂₈H₃₃N₃O₇).



Mol	Chain	Residues	Atoms				AltConf
57	A	1	Total	C	N	O	0
			38	28	3	7	

Chain 4:  37% 63%

MET ALA ASN LEU PHE ILE ARG LYS MET VAL ASN PRO LEU TYR SER ARG HIS THR VAL LYS PRO ARG ALA LEU SER THR PHE LEU PHE GLY SER ILE ARG GLY ALA VAL PRO VAL ALA VAL PRO GLY ALA VAL ARG SER LEU SER PRO GLY LEU PRO HIS LEU

LEU PRO ALA LEU GLY F66 M103


- Molecule 6: 39S ribosomal protein L37, mitochondrial

Chain 5:  6% 93% 7%

MET ALA LEU ALA SER GLY PRO ALA ARG ALA LEU ALA LEU GLY SER GLY GLN LEU LEU GLY LEU GLY PHE GLY ALA PRO ARG ARG GLY A30 D47 G54 L55 E56 G81 D101 G131 E134 L137 V140 D141 D142 H146 L145 D182 E243 Y257

G282 R300 E337 D365 C366 N367 E368 D377 S378 D379 K394 R395 A422 ALA


- Molecule 7: 39S ribosomal protein L38, mitochondrial

Chain 6:  12% 92% 7%

MET ALA ALA PRO TRP ARG ALA ALA LEU CYS CYS ARG ARG TRP ARG ALA PHE SER THR SER ALA VAL LEU GLY E27 D39 R47 R52 E62 E80 K81 T82 D83 P84 K85 E86 D89 K96 R99 A115 R121 A122 L125 R126 T127 D133

A178 V179 G180 E181 D182 E197 A208 E209 E210 D229 A230 E231 R244 V245 A246 E247 K274 Q275 D276 S282 E283 D284 A285 C290 R299 D302 K306 H307 D325 E339 H373 Y380

- Molecule 8: 39S ribosomal protein L39, mitochondrial

Chain 7:  9% 86% 14%

MET GLU ALA LEU MET GLY SER ARG ALA LEU ARG ALA TRP LEU VAL ALA PRO GLY GLY GLY ILE ARG PHE ILE ALA THR SER SER ALA SER GLN LEU S36 E51 L58 T59 P60 R61 T62 E63 E66 V67 K68 D114 C133 N165 D182 S187 K188

E191 K204 E220 E241 A244 S245 Q246 N247 R250 T257 Q258 D259 L286 T289 Q290 P291 G299 D314 E326 ASP GLN SER SER LYS ALA THR GLU ILE GLU PRO MET ARG SER THR

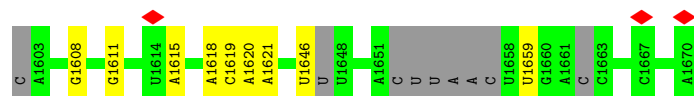
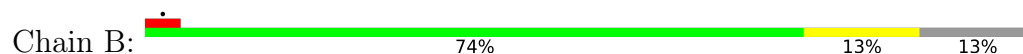
- Molecule 9: 39S ribosomal protein L40, mitochondrial

Chain 8:  17% 50% 49%

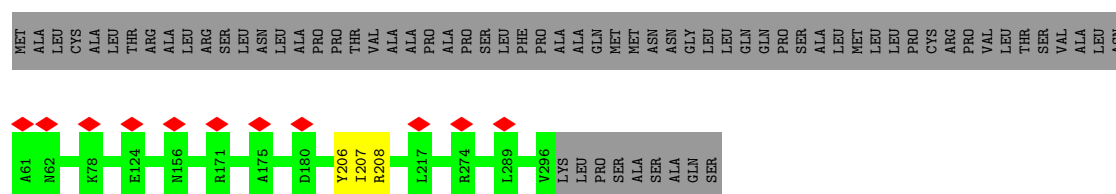
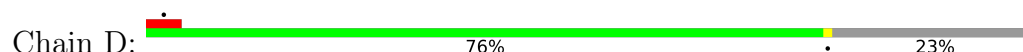
MET THR ALA SER VAL LEU ARG SER ILE SER SER LEU ALA ARG ARG PRO THR THR GLY LEU LEU GLY THR TRP GLN THR GLN ARG GLU THR HIS GLN ARG ALA SER LEU LEU SER PHE TRP LEU LEU ILE PRO MET ARG SER PRO ARG LEU ARG LYS LYS LYS VAL ASP PRO LYS LYS



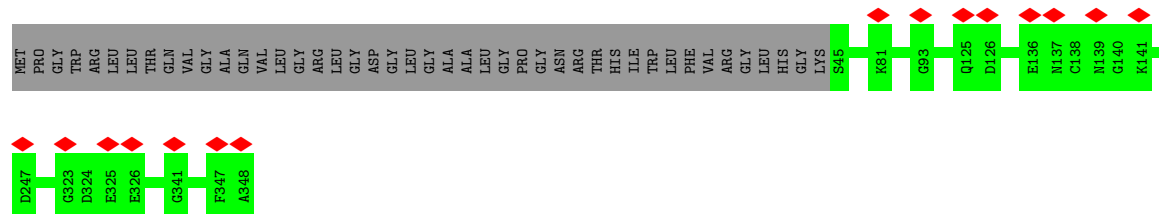
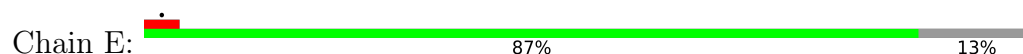
- Molecule 12: tRNA Val



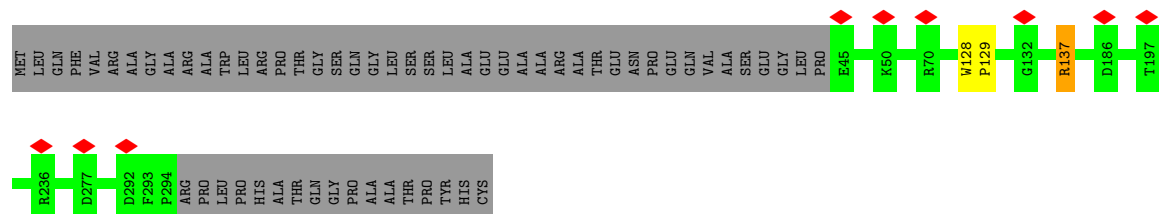
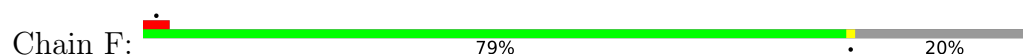
- Molecule 13: 39S ribosomal protein L2, mitochondrial



- Molecule 14: 39S ribosomal protein L3, mitochondrial

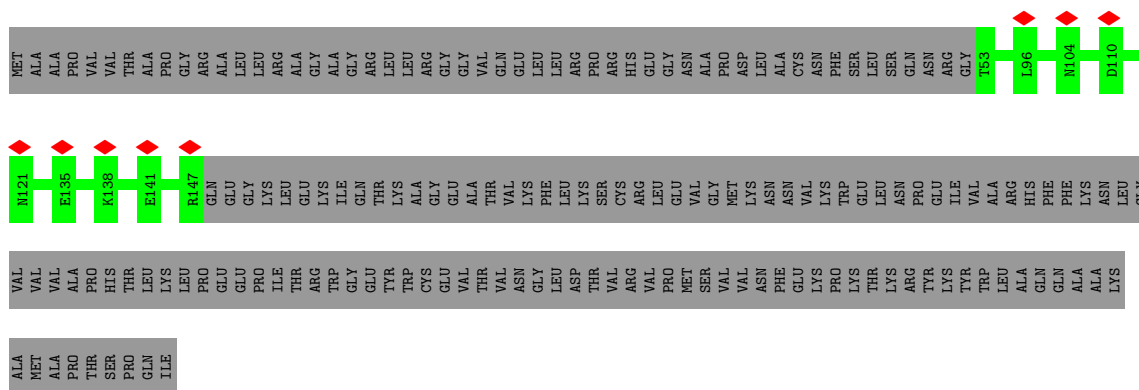


- Molecule 15: 39S ribosomal protein L4, mitochondrial

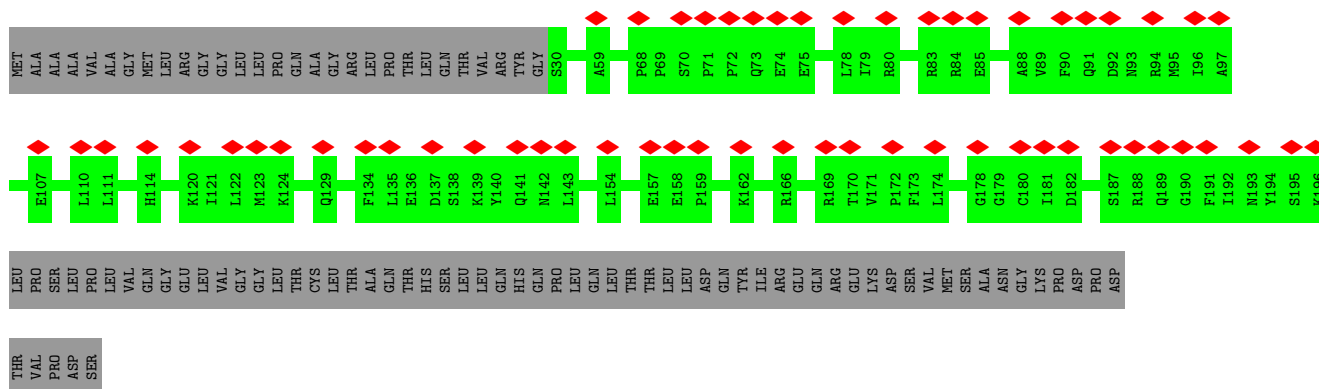


- Molecule 16: 39S ribosomal protein L9, mitochondrial

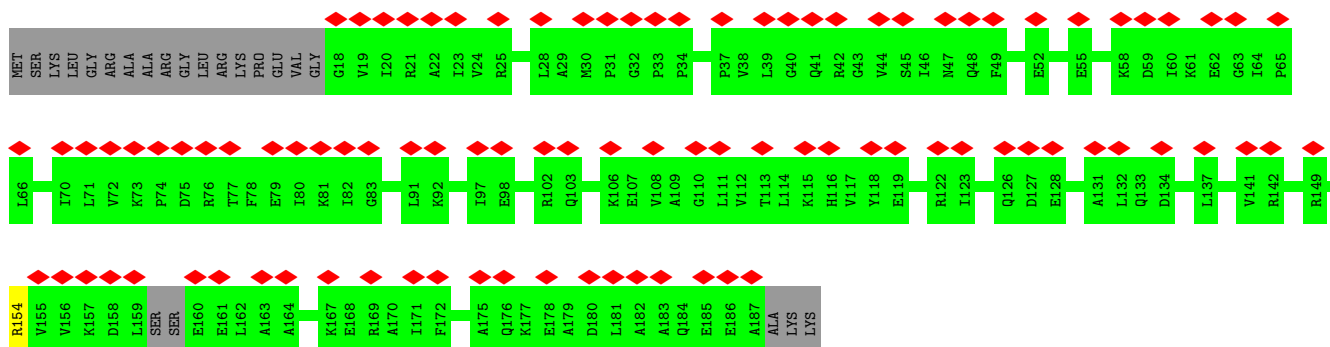
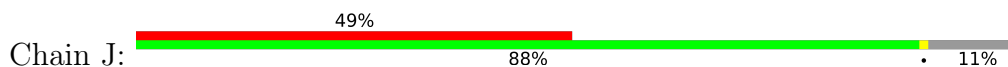




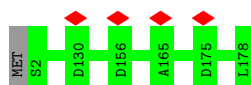
- Molecule 17: 39S ribosomal protein L10, mitochondrial



- Molecule 18: 39S ribosomal protein L11, mitochondrial

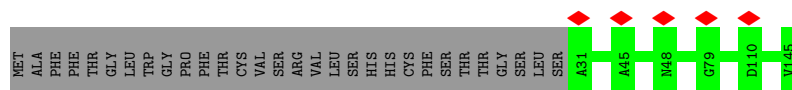


- Molecule 19: 39S ribosomal protein L13, mitochondrial



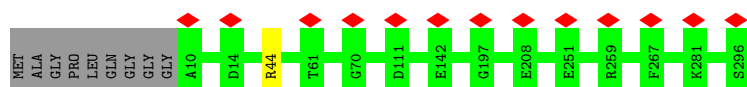
- Molecule 20: 39S ribosomal protein L14, mitochondrial

Chain L:  79% 21%




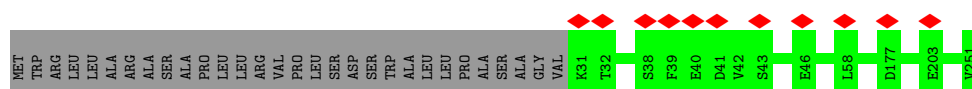
- Molecule 21: 39S ribosomal protein L15, mitochondrial

Chain M:  97% .



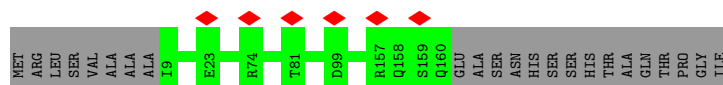
- Molecule 22: 39S ribosomal protein L16, mitochondrial

Chain N:  88% 12%




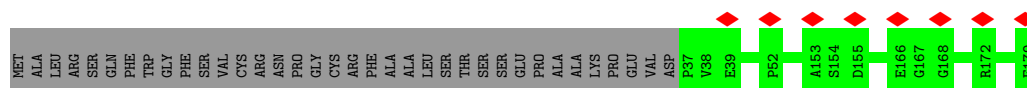
- Molecule 23: 39S ribosomal protein L17, mitochondrial

Chain O:  87% 13%




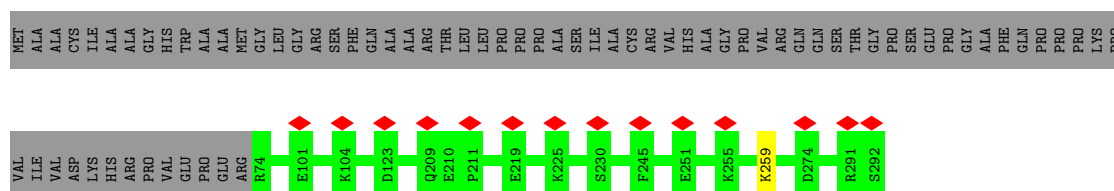
- Molecule 24: Mitochondrial ribosomal protein L18, isoform CRA_b

Chain P:  80% 20%



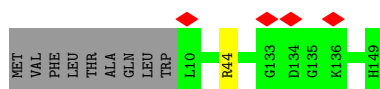
- Molecule 25: 39S ribosomal protein L19, mitochondrial

Chain Q:  5% 75% 25%




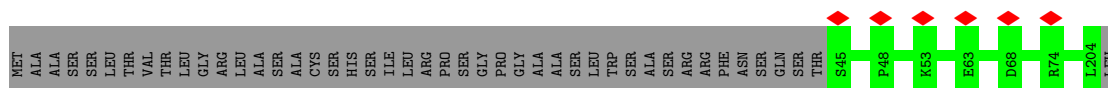
- Molecule 26: 39S ribosomal protein L20, mitochondrial

Chain R:  93% 6%




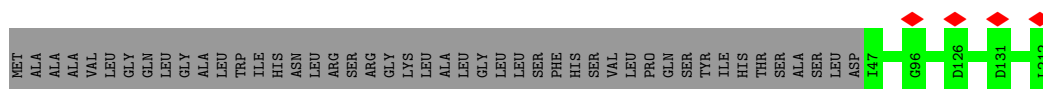
- Molecule 27: 39S ribosomal protein L21, mitochondrial

Chain S:  78% 22%

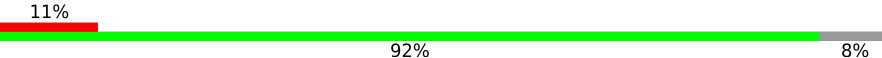


- Molecule 28: 39S ribosomal protein L22, mitochondrial

Chain T:  78% 22%



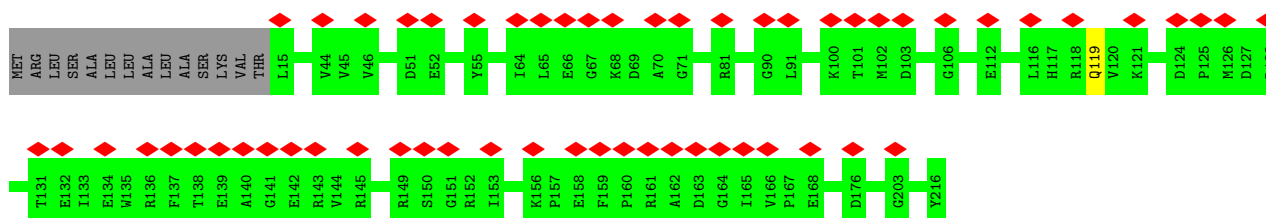
- Molecule 29: 39S ribosomal protein L23, mitochondrial

Chain U:  11% 92% 8%




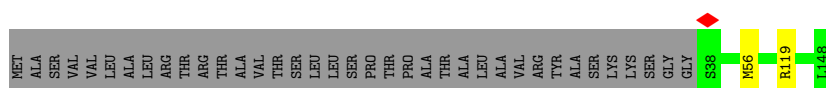
- Molecule 30: 39S ribosomal protein L24, mitochondrial

Chain V:  27% 93% 6%



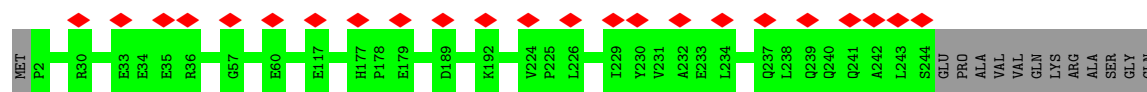
- Molecule 31: 39S ribosomal protein L27, mitochondrial

Chain W:  74% 25%

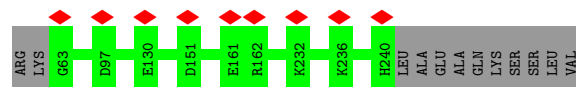
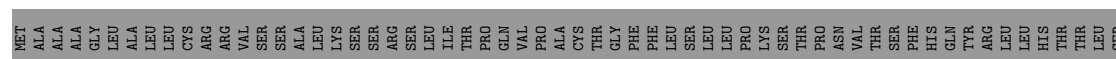


- Molecule 32: 39S ribosomal protein L28, mitochondrial

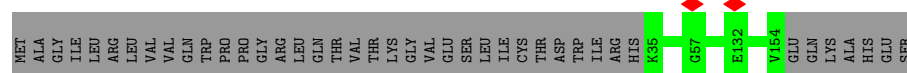
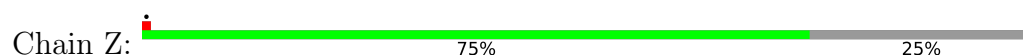
Chain X:  9% 95% 5%



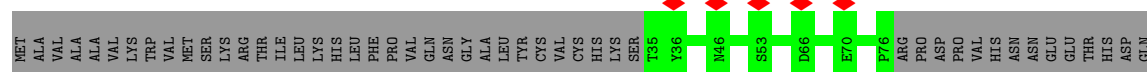
- Molecule 33: 39S ribosomal protein L47, mitochondrial



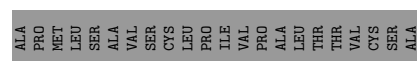
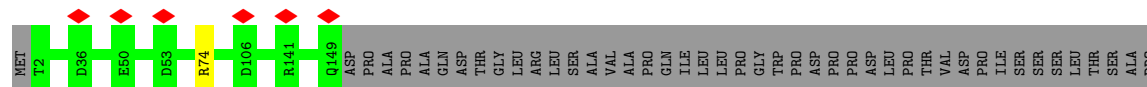
- Molecule 34: 39S ribosomal protein L30, mitochondrial



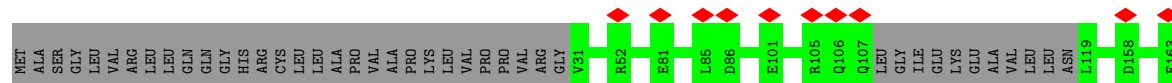
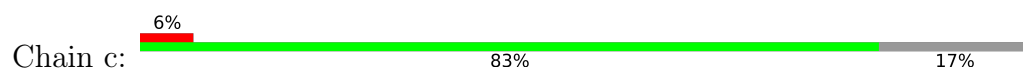
- Molecule 35: 39S ribosomal protein L42, mitochondrial



- Molecule 36: 39S ribosomal protein L43, mitochondrial

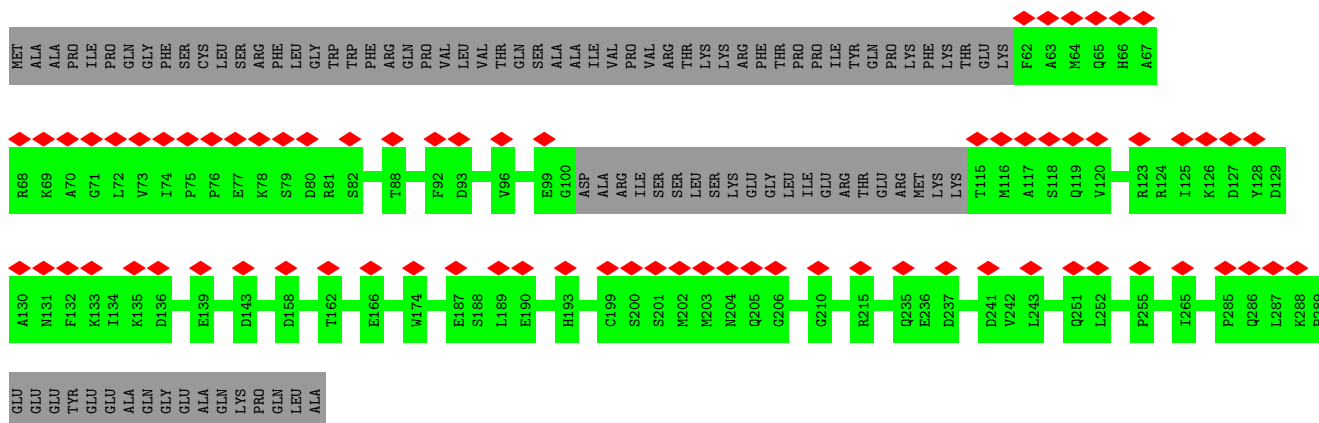


- Molecule 37: 39S ribosomal protein L44, mitochondrial

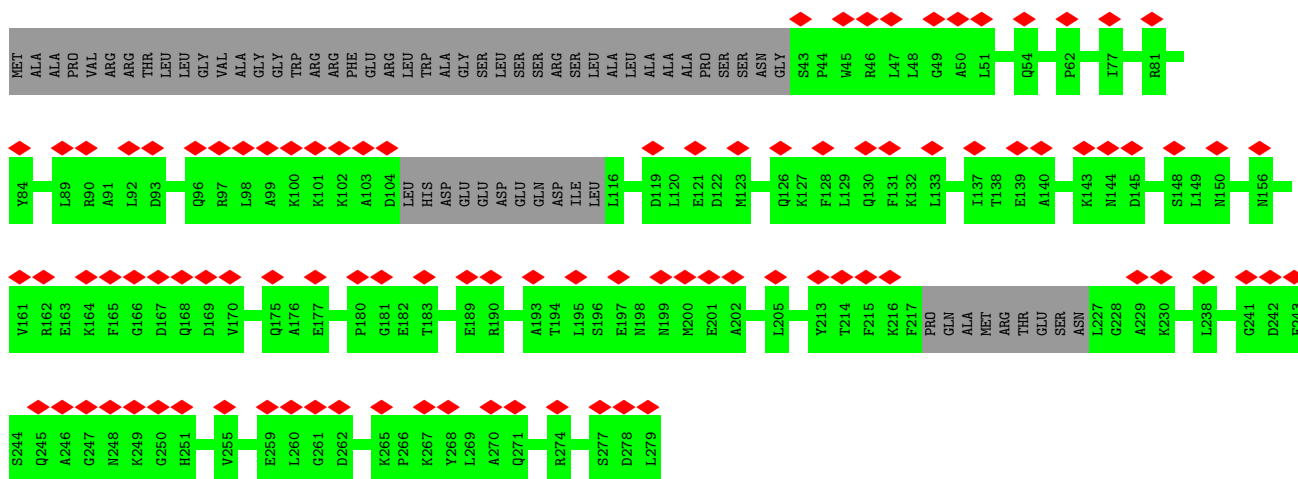
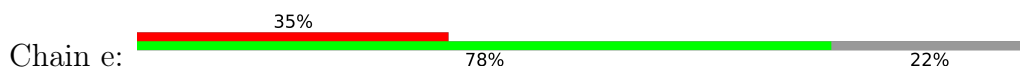




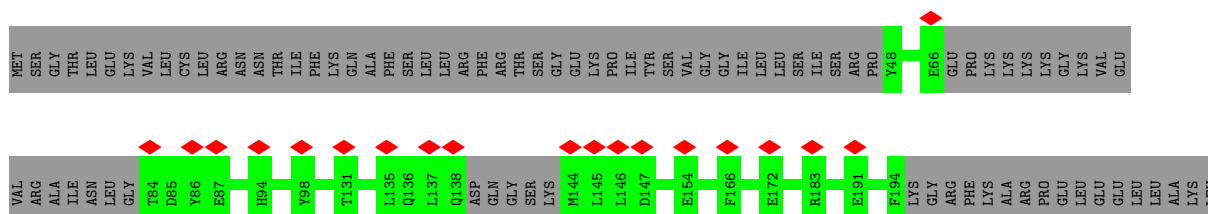
- Molecule 38: 39S ribosomal protein L45, mitochondrial



- Molecule 39: 39S ribosomal protein L46, mitochondrial

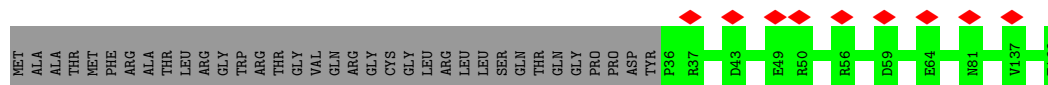
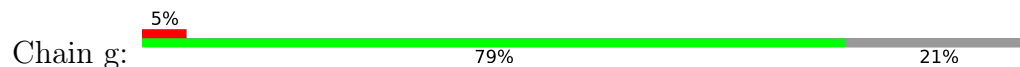


- Molecule 40: 39S ribosomal protein L48, mitochondrial

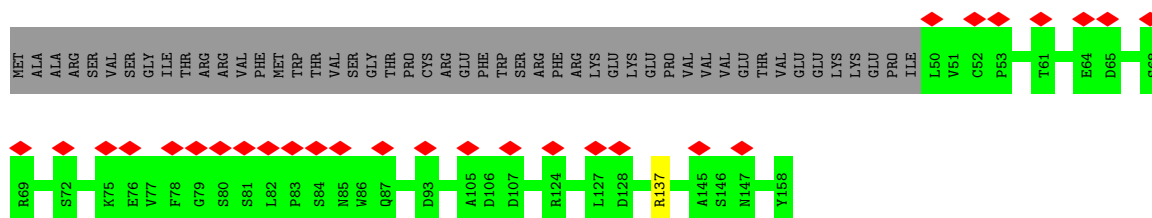


LYS

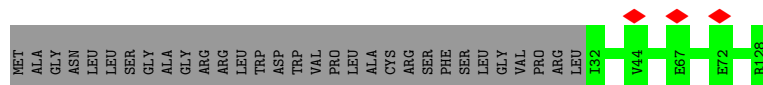
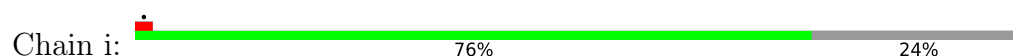
- Molecule 41: 39S ribosomal protein L49, mitochondrial



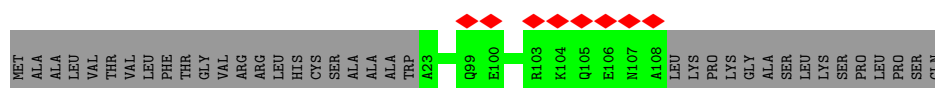
- Molecule 42: 39S ribosomal protein L50, mitochondrial



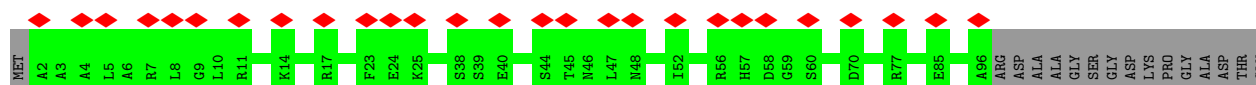
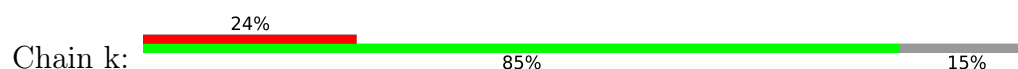
- Molecule 43: 39S ribosomal protein L51, mitochondrial



- Molecule 44: cDNA FLJ76418, highly similar to Homo sapiens mitochondrial ribosomal protein L52 (MRPL52), transcript variant 1, mRNA

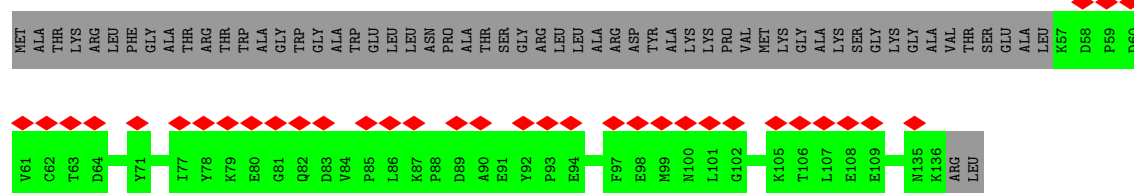


- Molecule 45: 39S ribosomal protein L53, mitochondrial

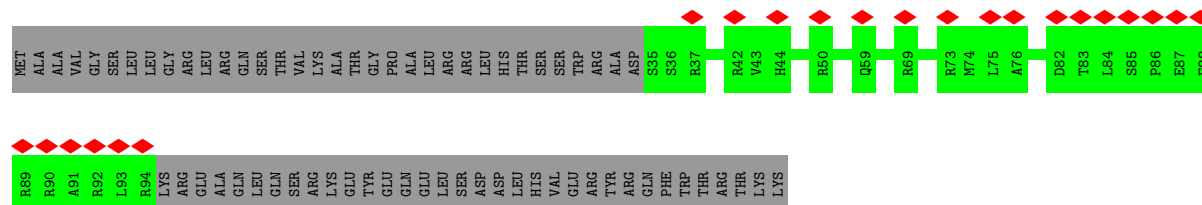


ARG

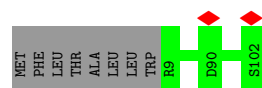
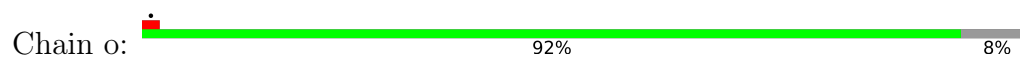
- Molecule 46: 39S ribosomal protein L54, mitochondrial



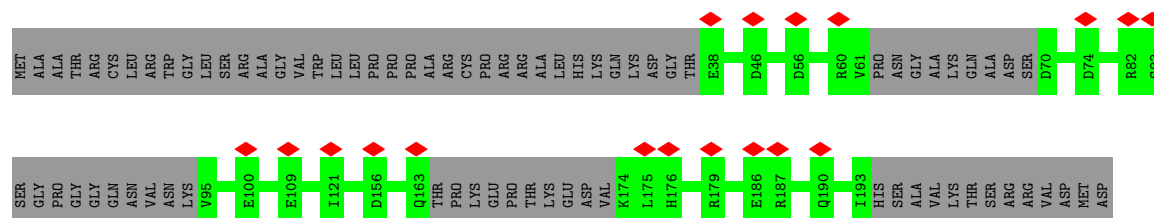
- Molecule 47: 39S ribosomal protein L55, mitochondrial



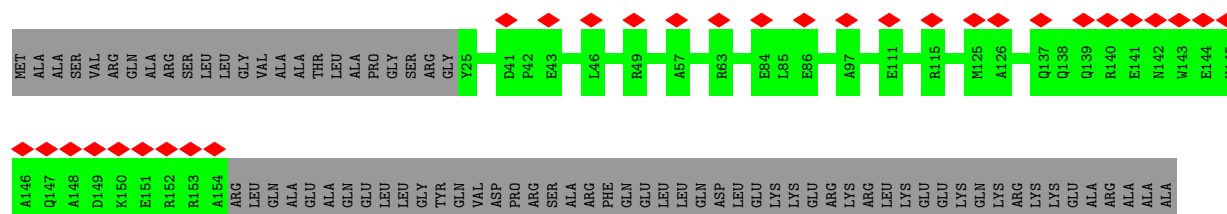
- Molecule 48: Ribosomal protein 63, mitochondrial



- Molecule 49: Peptidyl-tRNA hydrolase ICT1, mitochondrial

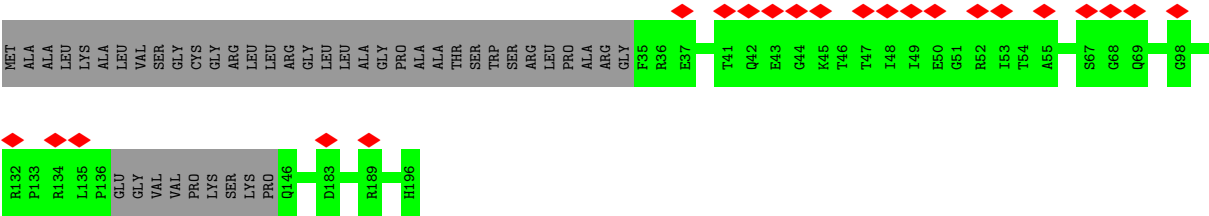
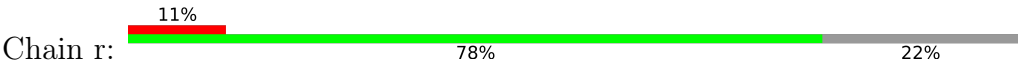


- Molecule 50: Growth arrest and DNA damage-inducible proteins-interacting protein 1

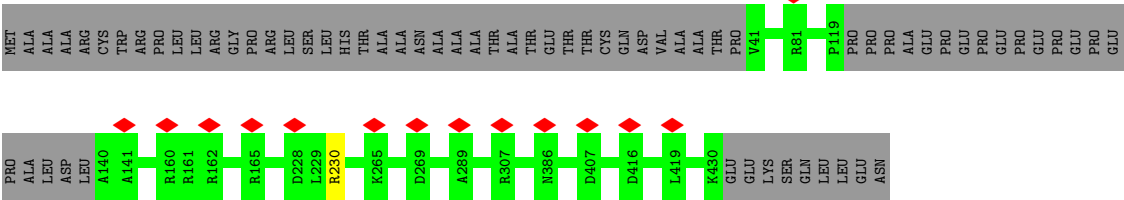
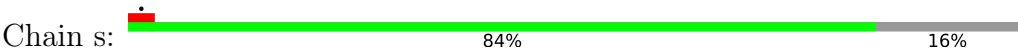


LEU
ALA
ALA
ALA
ALA
VAL
ALA
GLN
ASP
PRO
ALA
ALA
SER
GLY
ALA
PRO
SER
SER

- Molecule 51: 39S ribosomal protein S18a, mitochondrial



- Molecule 52: 39S ribosomal protein S30, mitochondrial



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	100000	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 ($e^-/\text{\AA}^2$)	38	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.166	Depositor
Minimum map value	-0.095	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	525.0, 525.0, 525.0	wwPDB
Map dimensions	500, 500, 500	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.05, 1.05, 1.05	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 5MU, ZN, 2MA, 2MU, H8Q, 5MC, H8T

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	0	0.25	0/895	0.55	0/1201
2	1	0.25	0/444	0.59	0/591
3	2	0.26	0/382	0.55	0/507
4	3	0.25	0/852	0.54	0/1136
5	4	0.25	0/348	0.59	0/458
6	5	0.25	0/3296	0.49	0/4486
7	6	0.26	0/3039	0.52	0/4131
8	7	0.26	0/2419	0.49	0/3267
9	8	0.27	0/909	0.49	0/1228
10	9	0.27	0/1024	0.49	0/1379
11	A	0.17	4/35562 (0.0%)	0.70	12/55341 (0.0%)
12	B	0.16	0/1422	0.69	0/2202
13	D	0.26	0/1879	0.56	0/2527
14	E	0.26	0/2464	0.49	0/3341
15	F	0.44	2/2071 (0.1%)	0.58	1/2817 (0.0%)
16	H	0.26	0/798	0.53	0/1073
17	I	0.26	0/1380	0.53	0/1864
18	J	0.27	0/1312	0.50	0/1768
19	K	0.26	0/1495	0.49	0/2029
20	L	0.25	0/904	0.53	0/1218
21	M	0.27	0/2359	0.55	0/3185
22	N	0.27	0/1825	0.53	0/2458
23	O	0.25	0/1269	0.54	0/1708
24	P	0.25	0/1189	0.55	0/1608
25	Q	0.26	0/1863	0.54	0/2509
26	R	0.25	0/1174	0.54	0/1572
27	S	0.26	0/1311	0.55	0/1778
28	T	0.26	0/1402	0.52	0/1886
29	U	0.26	0/1200	0.54	0/1623
30	V	0.27	0/1696	0.52	0/2299
31	W	0.25	0/893	0.50	0/1204
32	X	0.26	0/2081	0.49	0/2812

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	Y	0.26	0/1571	0.51	0/2106
34	Z	0.26	0/1003	0.48	0/1354
35	a	0.25	0/717	0.50	0/975
36	b	0.25	0/1202	0.53	0/1626
37	c	0.26	0/2264	0.48	0/3059
38	d	0.26	0/1783	0.50	0/2415
39	e	0.27	0/1797	0.51	0/2422
40	f	0.25	0/1016	0.46	0/1374
41	g	0.26	0/1121	0.52	0/1528
42	h	0.27	0/908	0.51	0/1235
43	i	0.25	0/849	0.53	0/1135
44	j	0.25	0/703	0.48	0/947
45	k	0.26	0/743	0.55	0/1003
46	l	0.26	0/694	0.46	0/942
47	m	0.26	0/507	0.60	0/682
48	o	0.26	0/818	0.54	0/1097
49	p	0.25	0/1071	0.53	0/1433
50	q	0.27	0/1123	0.53	0/1519
51	r	0.26	0/1291	0.52	0/1746
52	s	0.26	0/3114	0.51	0/4225
All	All	0.24	6/105452 (0.0%)	0.60	13/150029 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
13	D	0	1
15	F	0	1
All	All	0	2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	F	129	PRO	N-CA	13.83	1.70	1.47
11	A	2930	U	O3'-P	8.85	1.71	1.61
11	A	2987	U	O3'-P	7.99	1.70	1.61
11	A	2015	G	O3'-P	6.92	1.69	1.61
15	F	128	TRP	C-N	6.68	1.47	1.34

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	F	129	PRO	CA-N-CD	-10.05	97.43	111.50
11	A	2987	U	P-O3'-C3'	8.71	130.15	119.70
11	A	2297	A	P-O3'-C3'	8.11	129.44	119.70
11	A	2071	U	C2-N1-C1'	7.81	127.07	117.70
11	A	2720	A	P-O3'-C3'	6.57	127.59	119.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
13	D	206	TYR	Peptide
15	F	137	ARG	Sidechain

5.2 Too-close contacts [i](#)

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

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	0	106/188 (56%)	104 (98%)	2 (2%)	0	100	100
2	1	51/65 (78%)	51 (100%)	0	0	100	100
3	2	44/92 (48%)	43 (98%)	1 (2%)	0	100	100
4	3	93/188 (50%)	91 (98%)	2 (2%)	0	100	100
5	4	35/103 (34%)	33 (94%)	2 (6%)	0	100	100
6	5	385/423 (91%)	366 (95%)	19 (5%)	0	100	100
7	6	346/380 (91%)	326 (94%)	20 (6%)	0	100	100
8	7	283/338 (84%)	264 (93%)	19 (7%)	0	100	100
9	8	103/206 (50%)	98 (95%)	5 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	9	122/137 (89%)	116 (95%)	6 (5%)	0	100	100
13	D	234/305 (77%)	219 (94%)	13 (6%)	2 (1%)	17	54
14	E	300/348 (86%)	288 (96%)	12 (4%)	0	100	100
15	F	248/311 (80%)	233 (94%)	15 (6%)	0	100	100
16	H	93/267 (35%)	90 (97%)	3 (3%)	0	100	100
17	I	165/261 (63%)	156 (94%)	9 (6%)	0	100	100
18	J	168/192 (88%)	158 (94%)	10 (6%)	0	100	100
19	K	175/178 (98%)	169 (97%)	6 (3%)	0	100	100
20	L	113/145 (78%)	108 (96%)	5 (4%)	0	100	100
21	M	285/296 (96%)	266 (93%)	19 (7%)	0	100	100
22	N	219/251 (87%)	209 (95%)	10 (5%)	0	100	100
23	O	150/175 (86%)	150 (100%)	0	0	100	100
24	P	139/179 (78%)	133 (96%)	6 (4%)	0	100	100
25	Q	217/292 (74%)	211 (97%)	6 (3%)	0	100	100
26	R	138/149 (93%)	133 (96%)	5 (4%)	0	100	100
27	S	158/205 (77%)	155 (98%)	3 (2%)	0	100	100
28	T	164/212 (77%)	160 (98%)	4 (2%)	0	100	100
29	U	137/153 (90%)	133 (97%)	4 (3%)	0	100	100
30	V	198/216 (92%)	190 (96%)	8 (4%)	0	100	100
31	W	109/148 (74%)	105 (96%)	4 (4%)	0	100	100
32	X	241/256 (94%)	238 (99%)	3 (1%)	0	100	100
33	Y	176/250 (70%)	172 (98%)	4 (2%)	0	100	100
34	Z	118/161 (73%)	112 (95%)	6 (5%)	0	100	100
35	a	79/142 (56%)	76 (96%)	3 (4%)	0	100	100
36	b	146/215 (68%)	136 (93%)	10 (7%)	0	100	100
37	c	271/332 (82%)	262 (97%)	9 (3%)	0	100	100
38	d	206/306 (67%)	201 (98%)	5 (2%)	0	100	100
39	e	211/279 (76%)	205 (97%)	6 (3%)	0	100	100
40	f	119/212 (56%)	116 (98%)	3 (2%)	0	100	100
41	g	129/166 (78%)	125 (97%)	4 (3%)	0	100	100
42	h	105/158 (66%)	103 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
43	i	95/128 (74%)	91 (96%)	4 (4%)	0	100	100
44	j	84/123 (68%)	83 (99%)	1 (1%)	0	100	100
45	k	93/112 (83%)	91 (98%)	2 (2%)	0	100	100
46	l	78/138 (56%)	76 (97%)	2 (3%)	0	100	100
47	m	58/128 (45%)	54 (93%)	4 (7%)	0	100	100
48	o	92/102 (90%)	88 (96%)	4 (4%)	0	100	100
49	p	119/206 (58%)	118 (99%)	1 (1%)	0	100	100
50	q	128/222 (58%)	128 (100%)	0	0	100	100
51	r	145/196 (74%)	138 (95%)	7 (5%)	0	100	100
52	s	366/439 (83%)	356 (97%)	10 (3%)	0	100	100
All	All	8037/10674 (75%)	7727 (96%)	308 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
13	D	207	ILE
13	D	208	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	0	97/164 (59%)	97 (100%)	0	100	100
2	1	50/60 (83%)	50 (100%)	0	100	100
3	2	40/72 (56%)	40 (100%)	0	100	100
4	3	88/166 (53%)	88 (100%)	0	100	100
5	4	37/89 (42%)	37 (100%)	0	100	100
6	5	353/368 (96%)	353 (100%)	0	100	100
7	6	313/332 (94%)	310 (99%)	3 (1%)	76	86
8	7	267/303 (88%)	267 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	8	97/190 (51%)	96 (99%)	1 (1%)	76	86
10	9	104/112 (93%)	104 (100%)	0	100	100
13	D	190/245 (78%)	190 (100%)	0	100	100
14	E	259/290 (89%)	259 (100%)	0	100	100
15	F	217/262 (83%)	216 (100%)	1 (0%)	88	93
16	H	86/228 (38%)	86 (100%)	0	100	100
17	I	154/232 (66%)	154 (100%)	0	100	100
18	J	134/150 (89%)	133 (99%)	1 (1%)	84	90
19	K	155/156 (99%)	155 (100%)	0	100	100
20	L	98/124 (79%)	98 (100%)	0	100	100
21	M	245/249 (98%)	244 (100%)	1 (0%)	91	94
22	N	188/211 (89%)	188 (100%)	0	100	100
23	O	133/150 (89%)	133 (100%)	0	100	100
24	P	125/154 (81%)	125 (100%)	0	100	100
25	Q	201/256 (78%)	200 (100%)	1 (0%)	88	93
26	R	118/126 (94%)	117 (99%)	1 (1%)	81	89
27	S	145/180 (81%)	145 (100%)	0	100	100
28	T	146/182 (80%)	146 (100%)	0	100	100
29	U	126/135 (93%)	126 (100%)	0	100	100
30	V	180/191 (94%)	179 (99%)	1 (1%)	86	91
31	W	91/119 (76%)	89 (98%)	2 (2%)	52	71
32	X	217/227 (96%)	217 (100%)	0	100	100
33	Y	161/223 (72%)	161 (100%)	0	100	100
34	Z	111/147 (76%)	111 (100%)	0	100	100
35	a	79/133 (59%)	79 (100%)	0	100	100
36	b	130/186 (70%)	129 (99%)	1 (1%)	81	89
37	c	241/288 (84%)	240 (100%)	1 (0%)	91	94
38	d	193/273 (71%)	193 (100%)	0	100	100
39	e	188/236 (80%)	188 (100%)	0	100	100
40	f	113/188 (60%)	113 (100%)	0	100	100
41	g	121/148 (82%)	121 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
42	h	103/148 (70%)	102 (99%)	1 (1%)	76	86
43	i	86/110 (78%)	86 (100%)	0	100	100
44	j	68/97 (70%)	68 (100%)	0	100	100
45	k	80/90 (89%)	80 (100%)	0	100	100
46	l	75/116 (65%)	75 (100%)	0	100	100
47	m	54/113 (48%)	54 (100%)	0	100	100
48	o	80/87 (92%)	80 (100%)	0	100	100
49	p	117/181 (65%)	117 (100%)	0	100	100
50	q	111/178 (62%)	111 (100%)	0	100	100
51	r	139/169 (82%)	139 (100%)	0	100	100
52	s	326/381 (86%)	325 (100%)	1 (0%)	92	95
All	All	7230/9215 (78%)	7214 (100%)	16 (0%)	93	96

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
42	h	137	ARG
37	c	302	ARG
26	R	44	ARG
36	b	74	ARG
25	Q	259	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
41	g	141	ASN
40	f	108	GLN
21	M	102	GLN
39	e	150	ASN
19	K	9	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
11	A	1490/1557 (95%)	252 (16%)	3 (0%)
12	B	55/69 (79%)	9 (16%)	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	1545/1626 (95%)	261 (16%)	3 (0%)

5 of 261 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
11	A	1685	C
11	A	1689	C
11	A	1692	A
11	A	1693	C
11	A	1695	C

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
11	A	2066	C
11	A	2558	A
11	A	2574	G

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

4 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	$\# Z > 2$
11	2MA	A	2990	11,54	17,25,26	1.09	1 (5%)	17,37,40	1.31	3 (17%)
11	5MC	A	2625	11	18,22,23	3.61	7 (38%)	26,32,35	0.99	2 (7%)
11	2MU	A	3039	11	20,23,24	8.02	15 (75%)	28,33,36	3.13	10 (35%)
11	5MU	A	2602	11	19,22,23	4.97	7 (36%)	28,32,35	3.62	10 (35%)

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
11	2MA	A	2990	11,54	-	2/3/25/26	0/3/3/3
11	5MC	A	2625	11	-	0/7/25/26	0/2/2/2
11	2MU	A	3039	11	-	0/9/27/28	0/2/2/2
11	5MU	A	2602	11	-	0/7/25/26	0/2/2/2

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	A	3039	2MU	C4-C5	21.83	1.81	1.44
11	A	3039	2MU	C6-N1	16.76	1.66	1.38
11	A	2602	5MU	C6-N1	11.86	1.58	1.38
11	A	3039	2MU	C6-C5	-11.78	1.15	1.34
11	A	3039	2MU	C4-N3	-10.56	1.19	1.38

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	A	2602	5MU	C5-C4-N3	12.45	125.94	115.31
11	A	3039	2MU	C5-C4-N3	9.41	123.34	115.31
11	A	2602	5MU	C5-C6-N1	-9.03	114.05	123.34
11	A	3039	2MU	C5-C6-N1	-7.04	116.09	123.34
11	A	3039	2MU	C4-N3-C2	-5.97	119.62	127.35

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	A	2990	2MA	O4'-C4'-C5'-O5'
11	A	2990	2MA	C3'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 140 ligands modelled in this entry, 137 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
56	H8Q	A	3430	-	77,80,80	1.11	4 (5%)	103,115,115	1.53	16 (15%)
57	H8T	A	3431	-	34,40,40	5.62	20 (58%)	37,55,55	2.64	10 (27%)
55	G	A	3429	-	18,25,26	1.03	2 (11%)	19,37,40	0.72	1 (5%)

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
56	H8Q	A	3430	-	-	33/83/127/127	0/8/8/8
57	H8T	A	3431	-	-	16/38/48/48	0/2/3/3
55	G	A	3429	-	-	0/3/25/26	0/3/3/3

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
57	A	3431	H8T	C4-N5	17.54	1.55	1.38
57	A	3431	H8T	C28-C29	12.67	1.62	1.32
57	A	3431	H8T	C4-C3	10.16	1.60	1.38
57	A	3431	H8T	C22-C23	9.16	1.55	1.32
57	A	3431	H8T	O36-C37	8.38	1.51	1.34

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
57	A	3431	H8T	C3-C2-C1	9.31	112.17	106.80
57	A	3431	H8T	O36-C37-C1	5.65	120.17	111.18
56	A	3430	H8Q	C3-C2-N71	-5.33	104.03	112.69
57	A	3431	H8T	C24-N25-C26	-5.06	113.85	122.03
56	A	3430	H8Q	C17-C18-N13	-4.85	105.17	109.92

There are no chirality outliers.

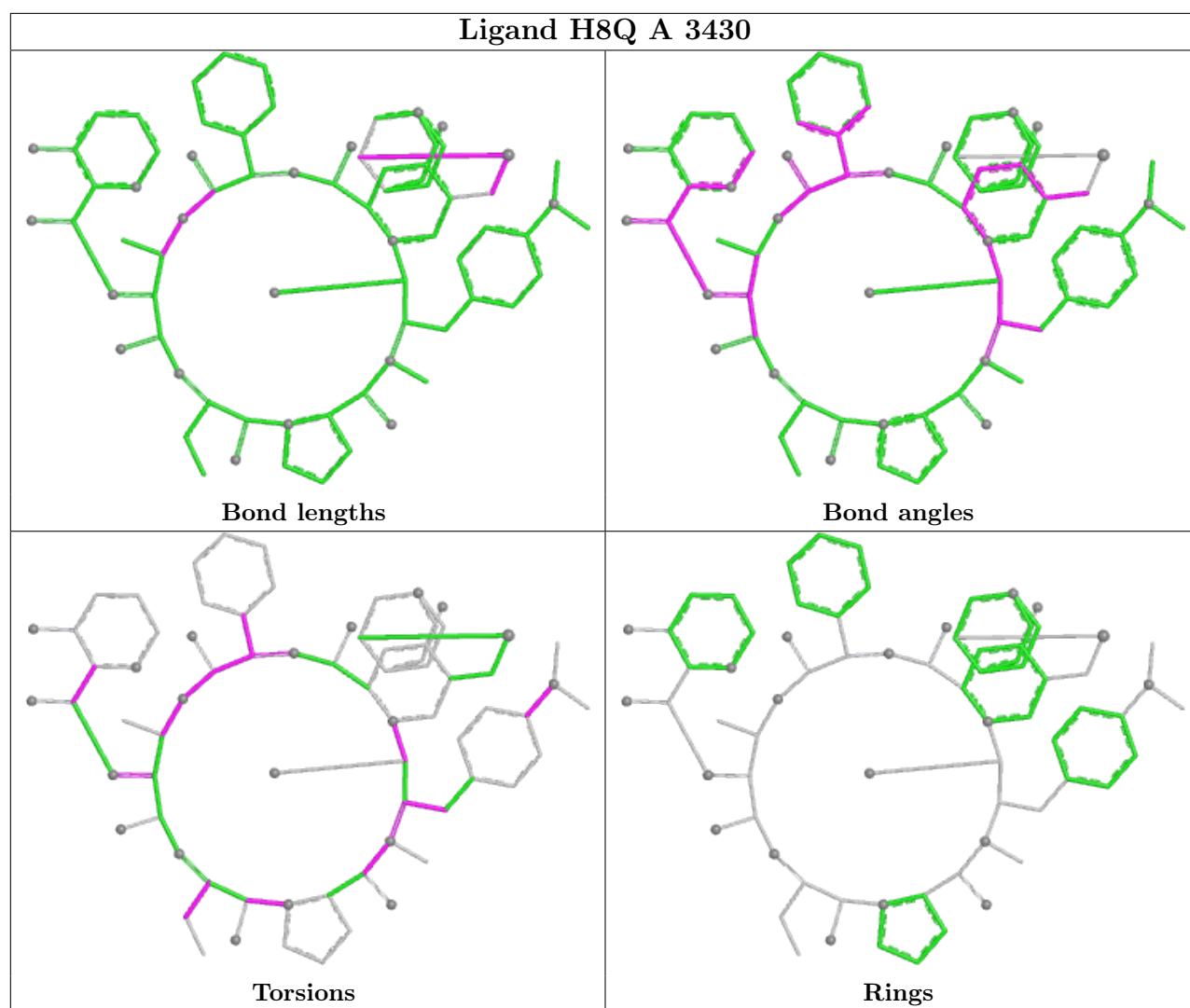
5 of 49 torsion outliers are listed below:

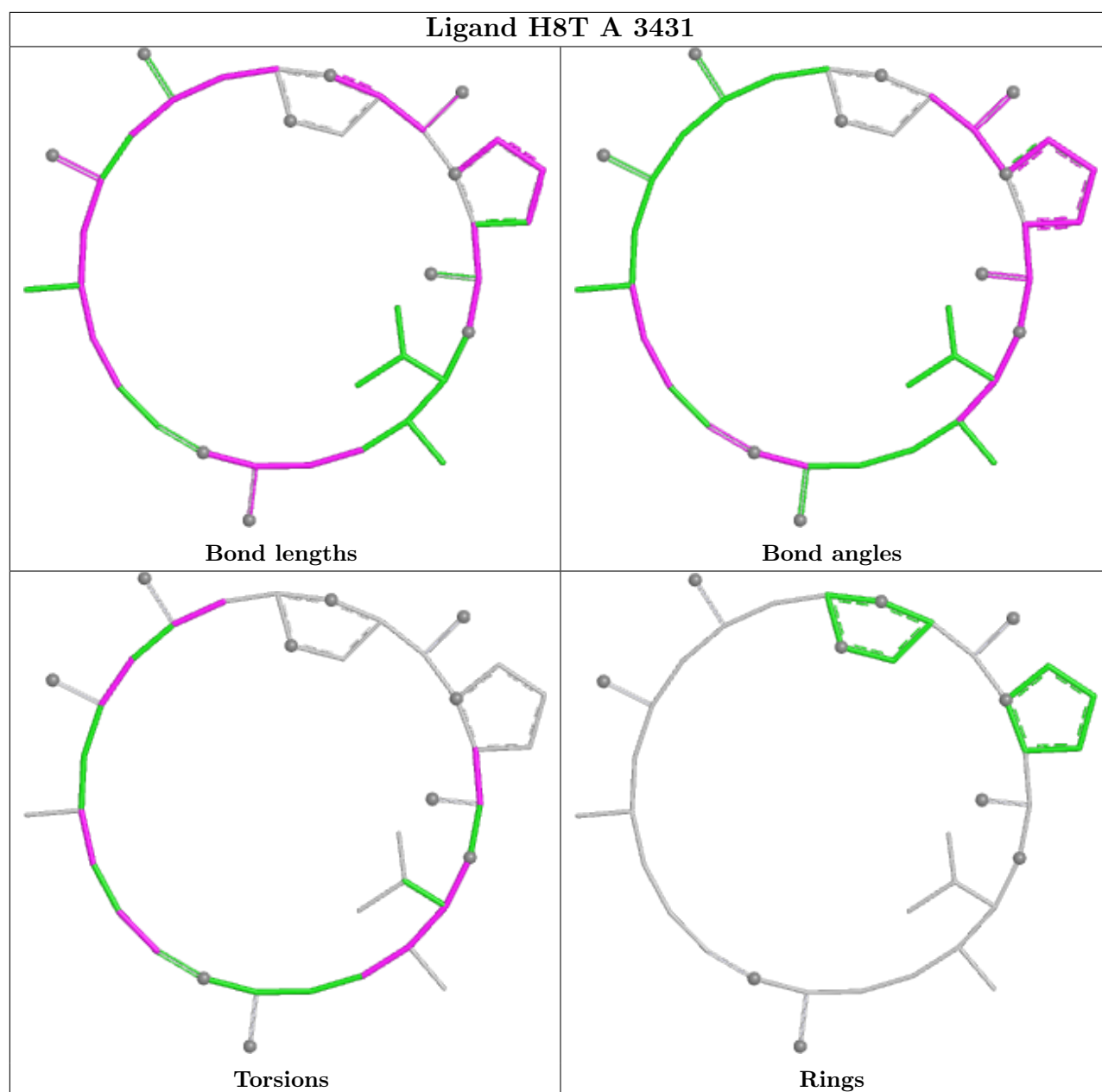
Mol	Chain	Res	Type	Atoms
56	A	3430	H8Q	C2-C1-N13-C14
56	A	3430	H8Q	O1-C1-N13-C14
56	A	3430	H8Q	O1-C1-N13-C18
56	A	3430	H8Q	C33-C34-O41-C43
56	A	3430	H8Q	O42-C34-O41-C43

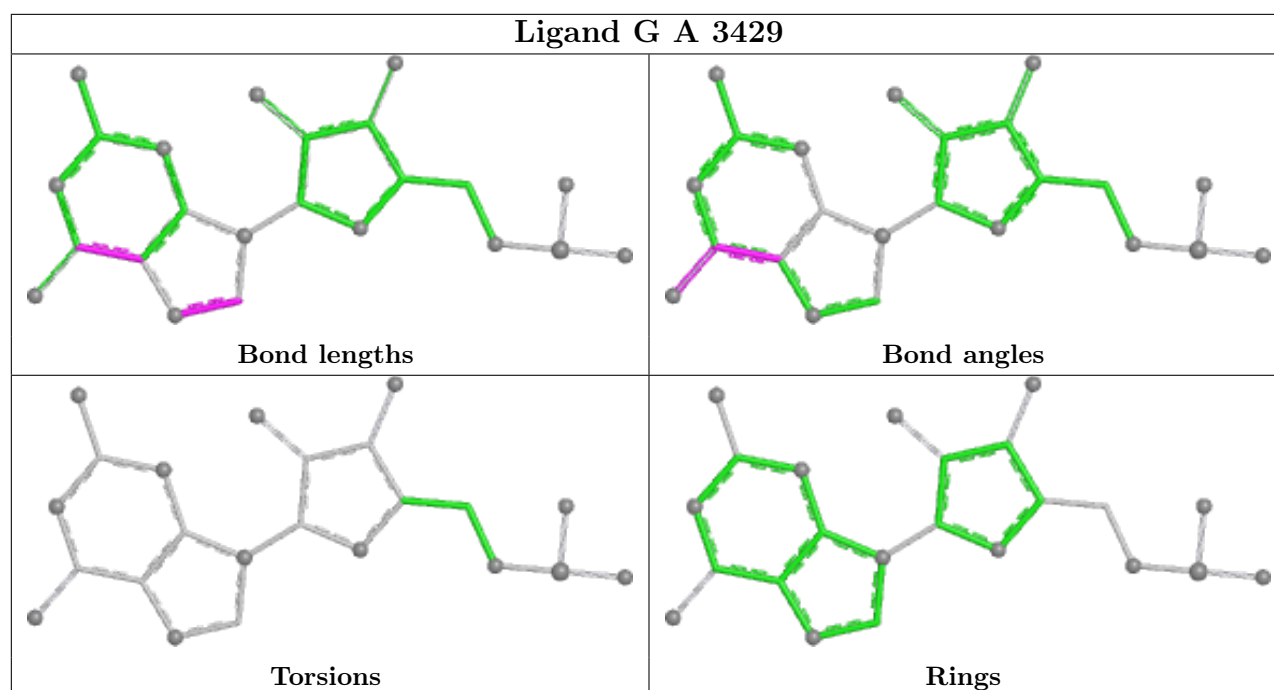
There are no ring outliers.

No monomer is involved in short contacts.

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
8	7	3
6	5	3
7	6	3
51	r	2
38	d	2
14	E	1
5	4	1
12	B	1
42	h	1
24	P	1
30	V	1
11	A	1

The worst 5 of 20 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	7	285:ASN	C	286:LEU	N	6.23
1	7	158:PHE	C	159:LYS	N	5.66
1	r	134:ARG	C	135:LEU	N	5.51
1	E	135:LYS	C	136:GLU	N	4.66
1	5	365:ASP	C	366:CYS	N	4.59

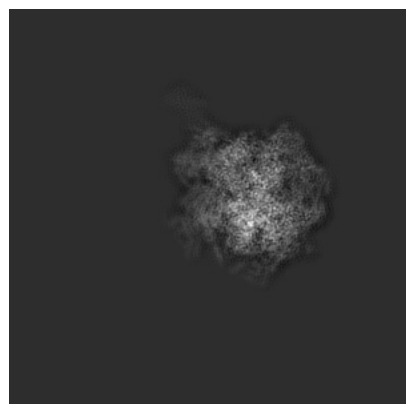
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4434. These allow visual inspection of the internal detail of the map and identification of artifacts.

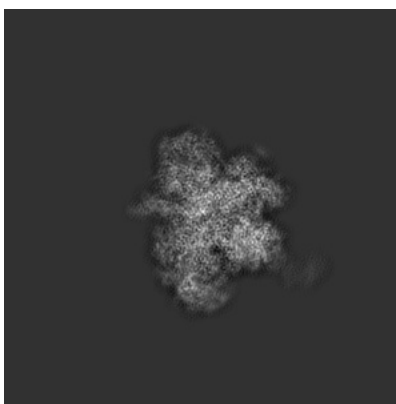
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

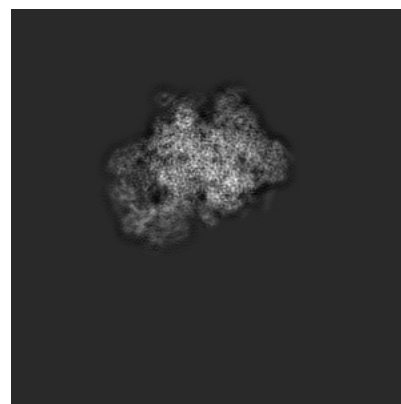
6.1.1 Primary map



X

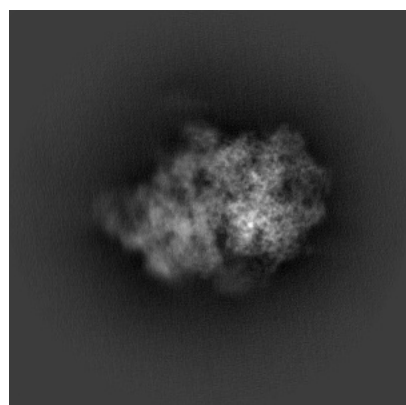


Y

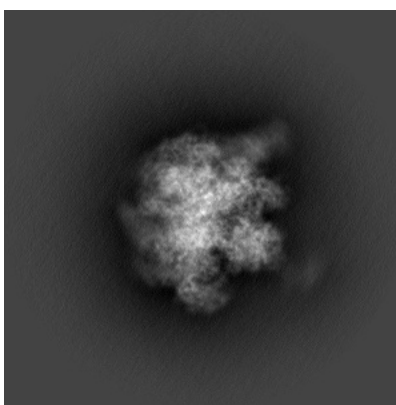


Z

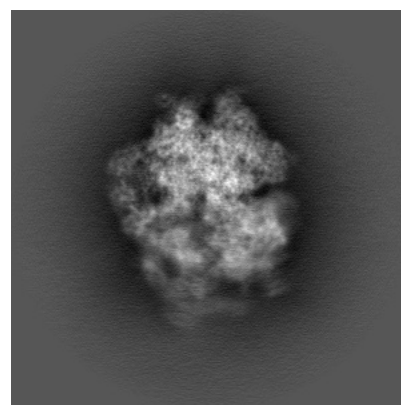
6.1.2 Raw map



X



Y

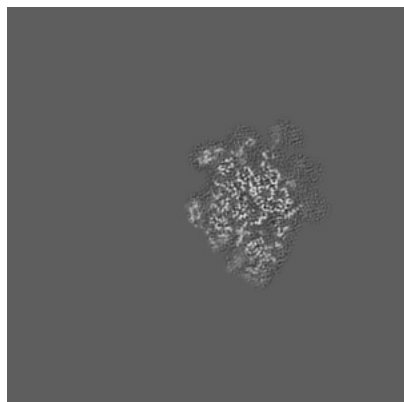


Z

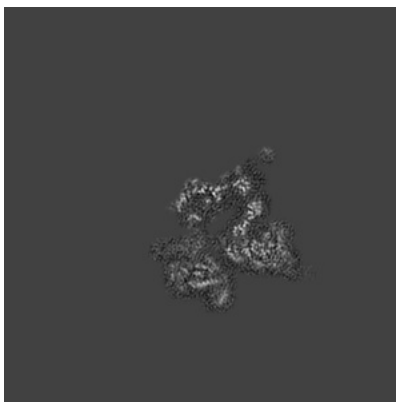
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

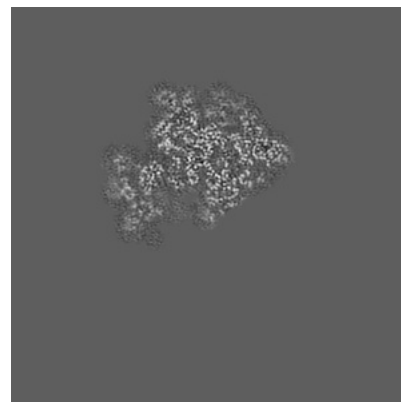
6.2.1 Primary map



X Index: 250

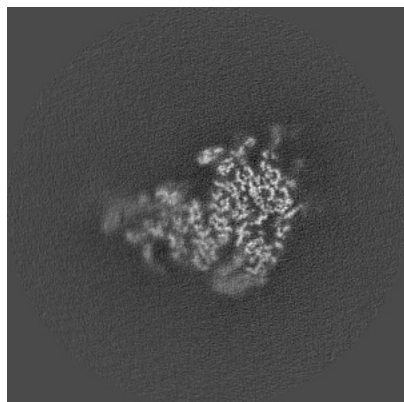


Y Index: 250

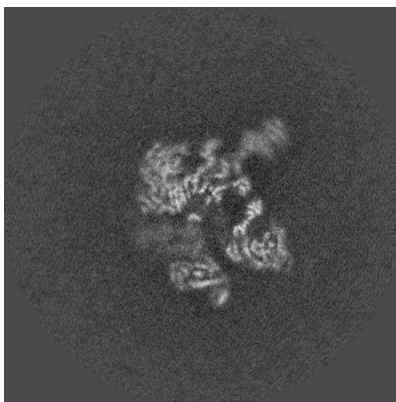


Z Index: 250

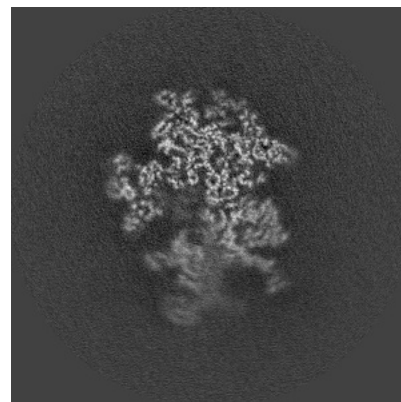
6.2.2 Raw map



X Index: 250



Y Index: 250

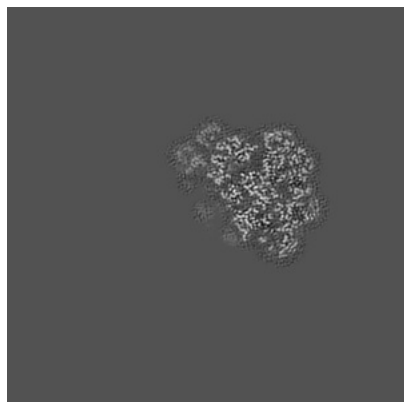


Z Index: 250

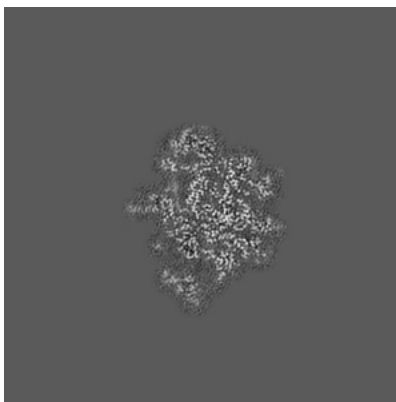
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

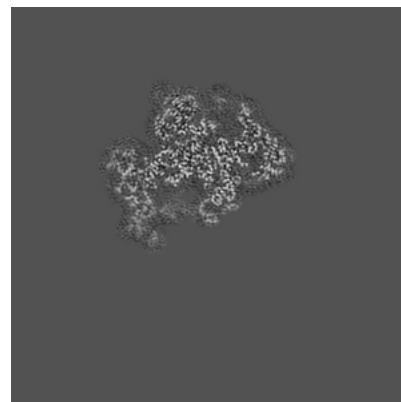
6.3.1 Primary map



X Index: 214

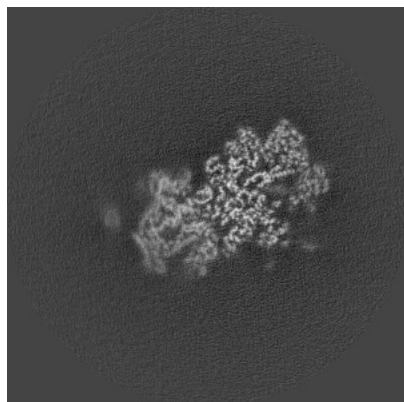


Y Index: 318

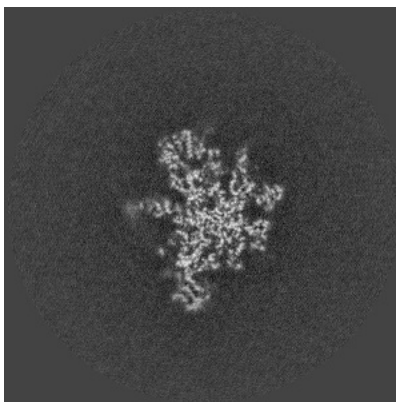


Z Index: 239

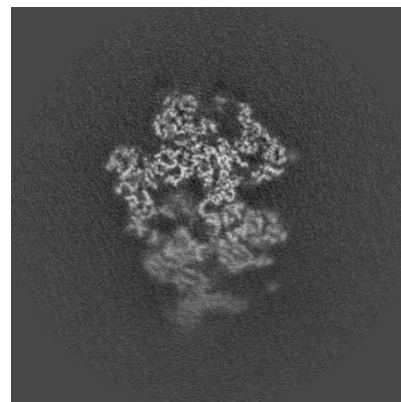
6.3.2 Raw map



X Index: 273



Y Index: 303

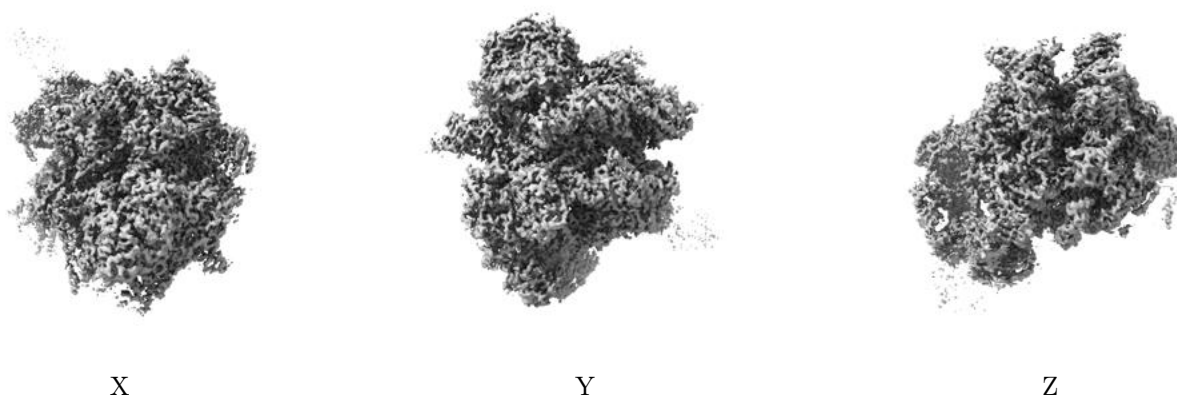


Z Index: 238

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

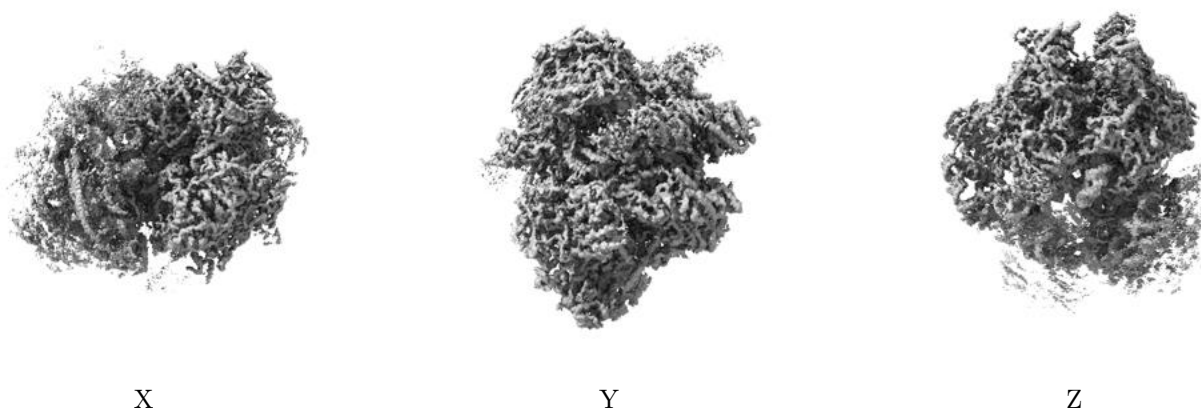
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

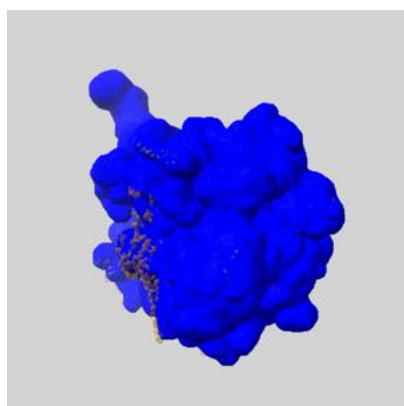
6.5 Mask visualisation [i](#)

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

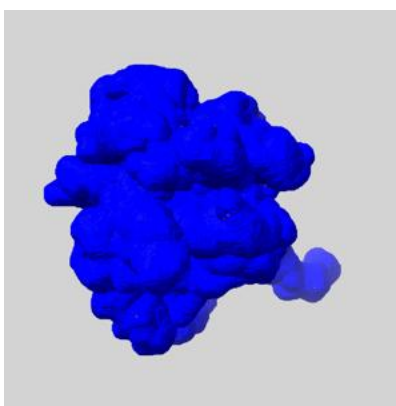
A mask typically either:

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

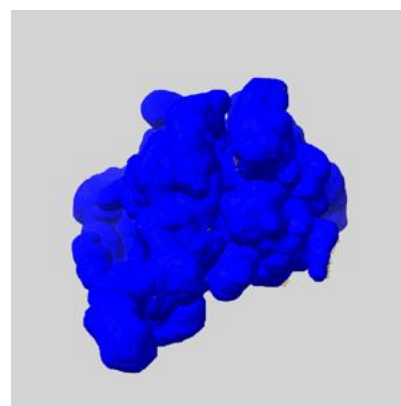
6.5.1 emd_4434_msk_1.map [i](#)



X



Y

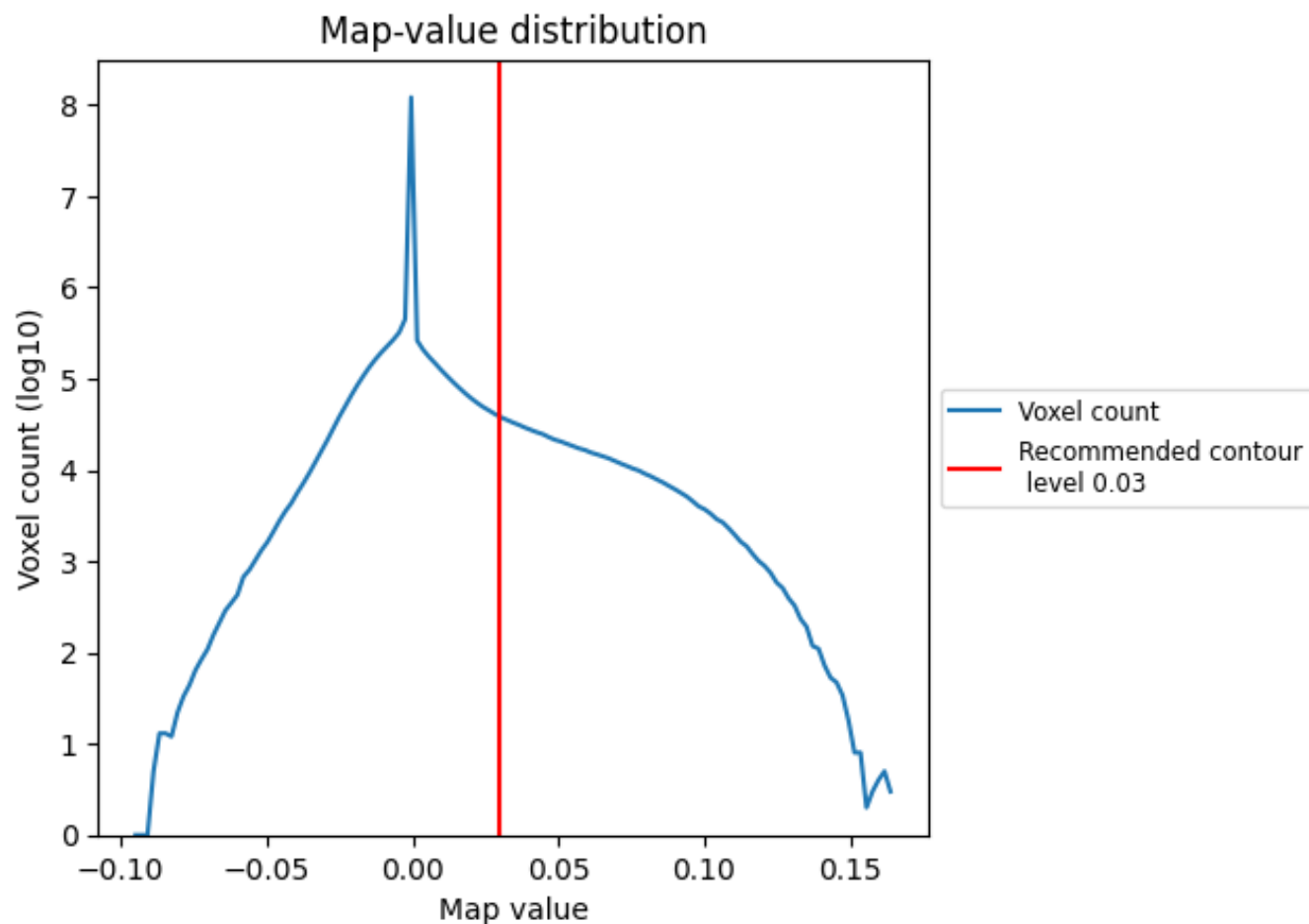


Z

7 Map analysis [i](#)

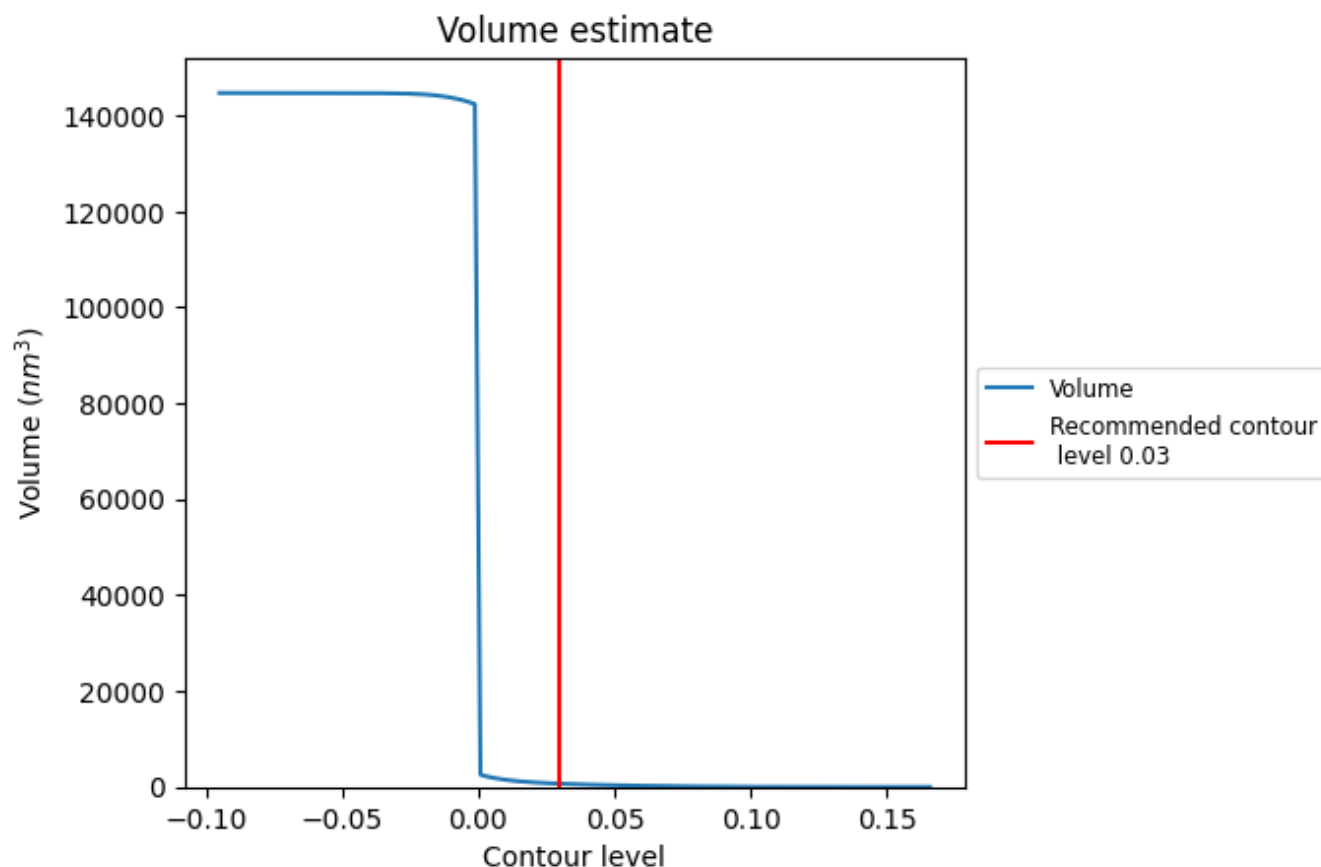
This section contains the results of statistical analysis of the map.

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.

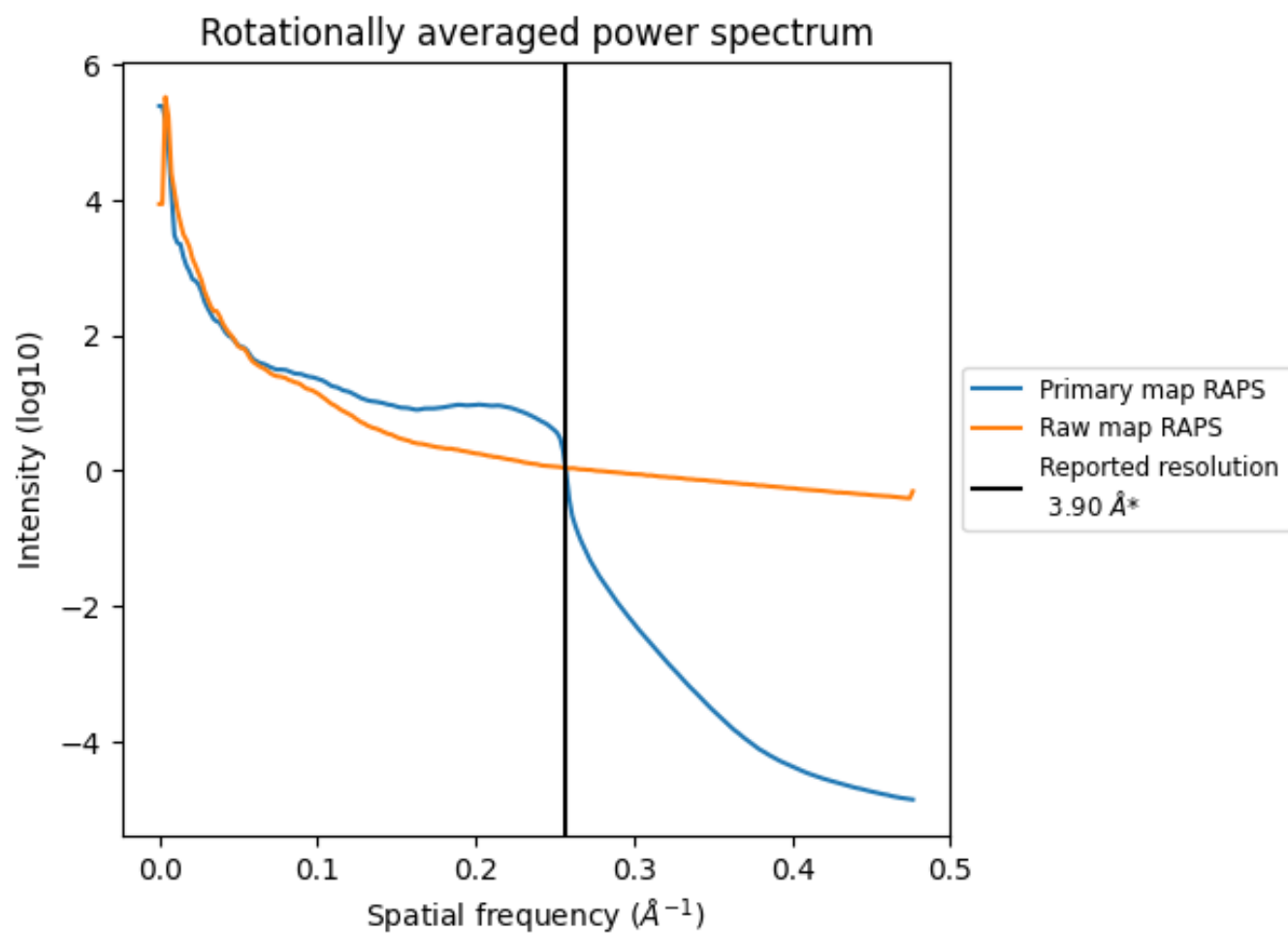
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 684 nm³; this corresponds to an approximate mass of 618 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 ⓘ

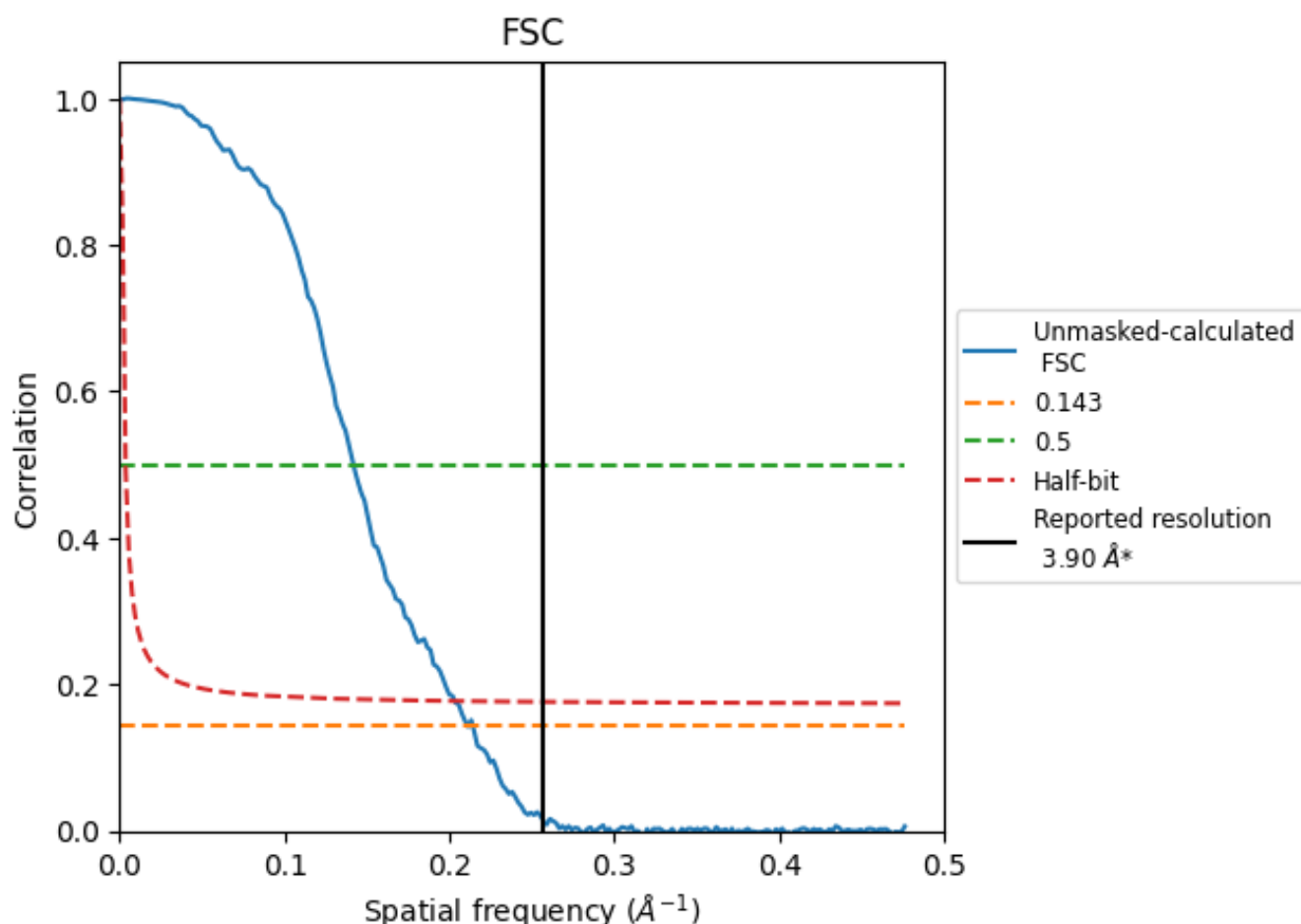


*Reported resolution corresponds to spatial frequency of 0.256 \AA^{-1}

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 0.256 Å⁻¹

8.2 Resolution estimates [i](#)

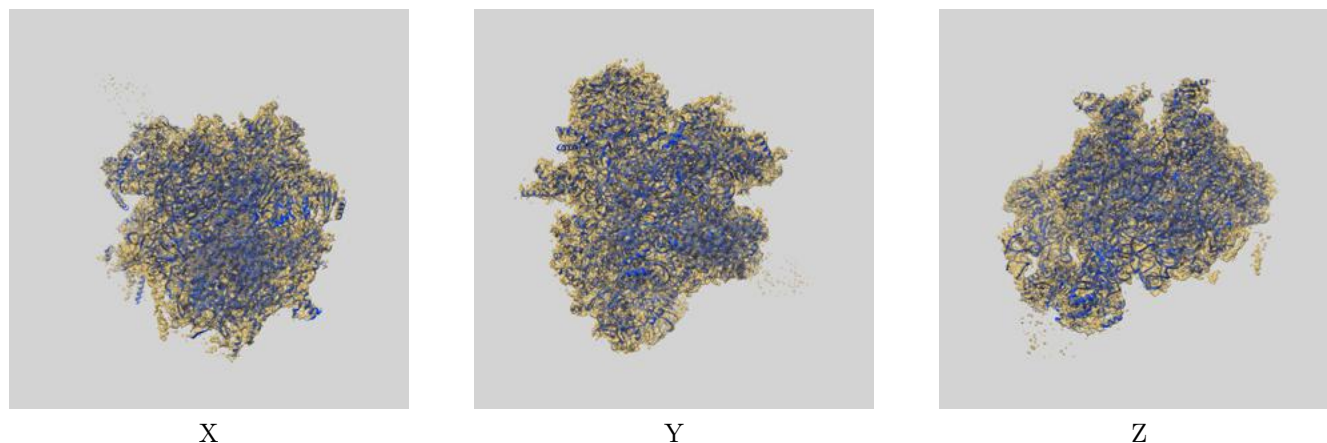
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.73	7.05	4.92

*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 4.73 differs from the reported value 3.9 by more than 10 %

9 Map-model fit [i](#)

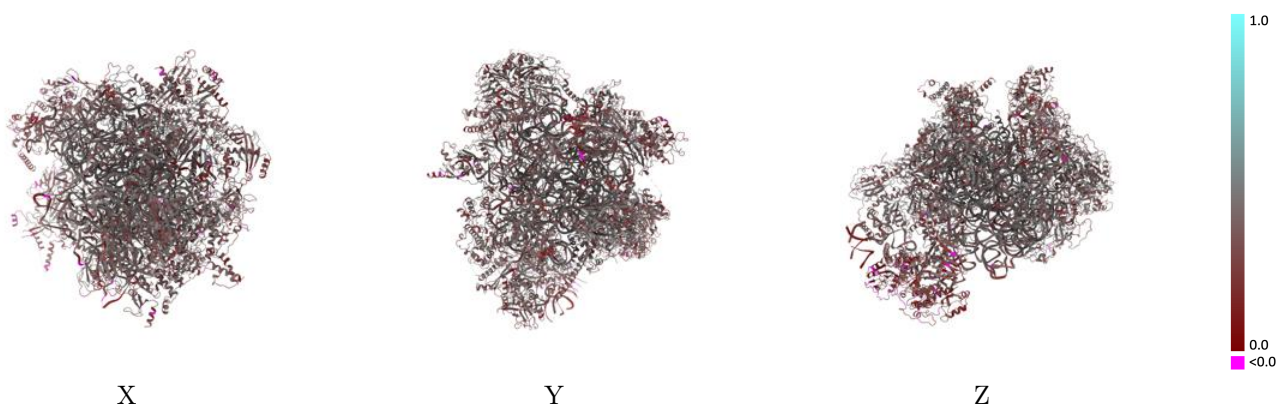
This section contains information regarding the fit between EMDB map EMD-4434 and PDB model 6I9R. Per-residue inclusion information can be found in section 3 on page 16.

9.1 Map-model overlay [i](#)



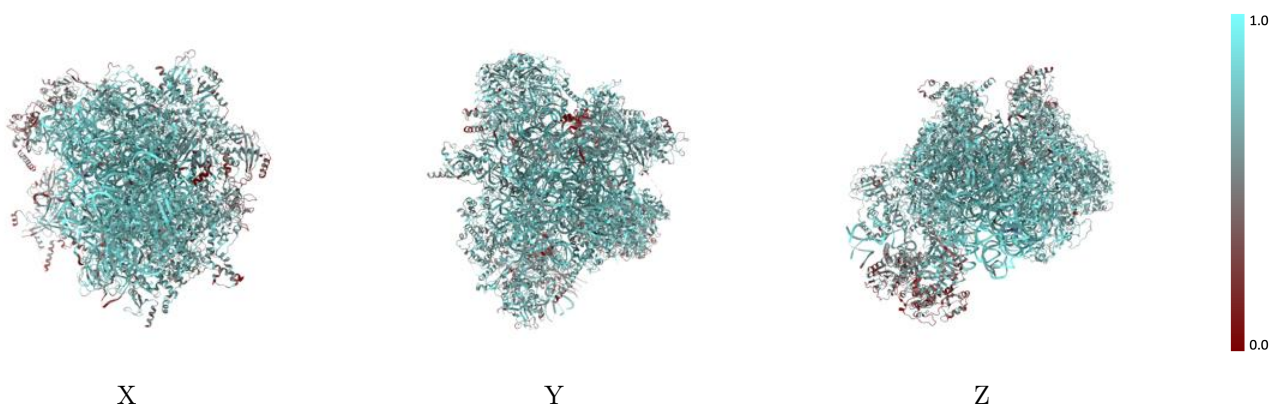
The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



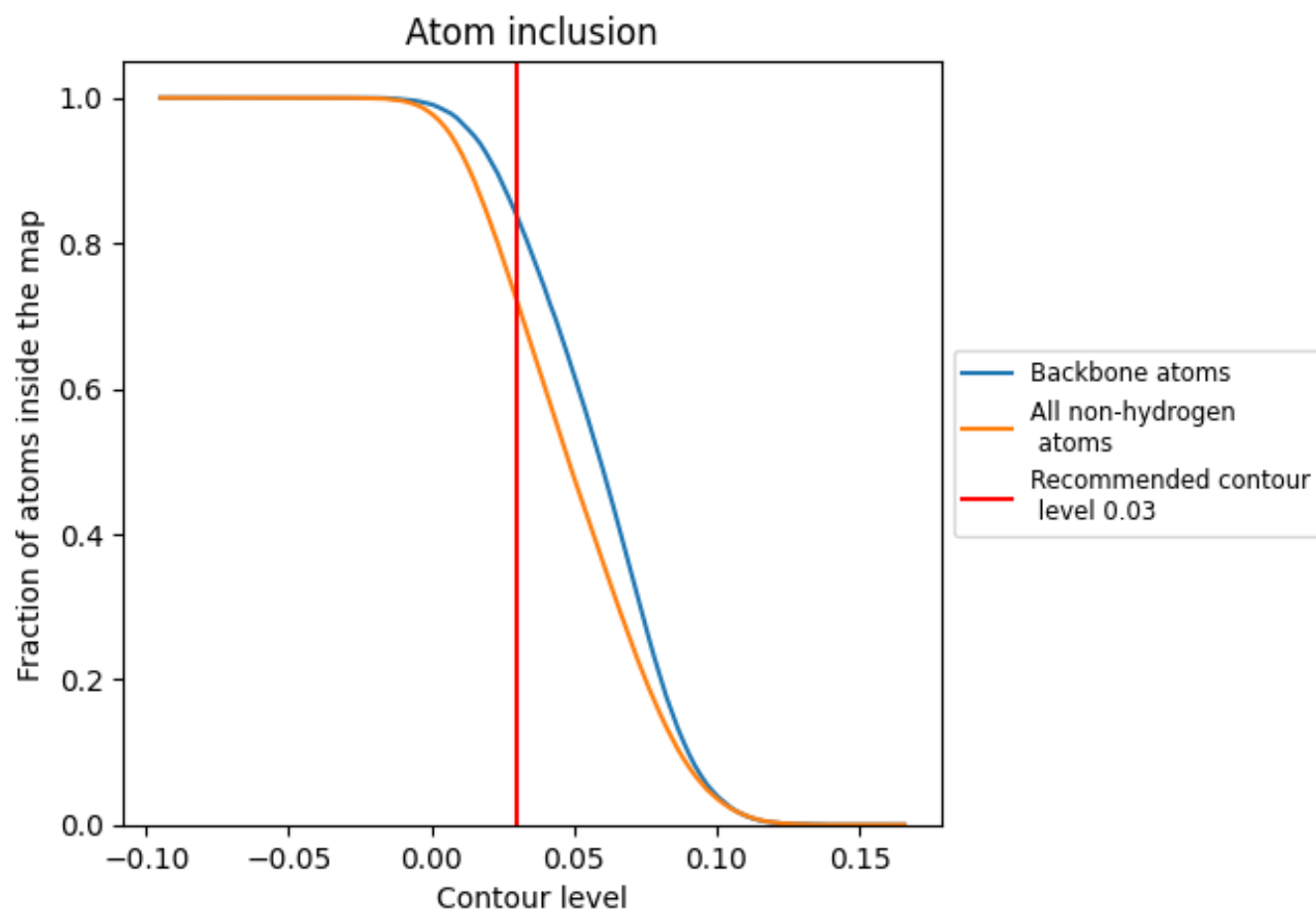
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.03).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ





































The table lists the average atom inclusion at the recommended contour level (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7206	 0.3740
0	 0.7022	 0.3720
1	 0.6573	 0.3550
2	 0.7472	 0.3870
3	 0.7469	 0.4220
4	 0.7638	 0.4080
5	 0.7045	 0.3770
6	 0.6625	 0.3390
7	 0.6550	 0.3480
8	 0.5012	 0.2580
9	 0.6300	 0.3660
A	 0.8442	 0.4030
B	 0.7953	 0.2910
D	 0.7293	 0.4020
E	 0.7203	 0.4010
F	 0.7157	 0.3950
H	 0.6398	 0.3610
I	 0.4977	 0.2820
J	 0.3425	 0.2230
K	 0.7403	 0.4100
L	 0.7491	 0.4170
M	 0.6958	 0.3790
N	 0.7187	 0.4020
O	 0.7250	 0.3860
P	 0.6889	 0.3610
Q	 0.6983	 0.3710
R	 0.7238	 0.3970
S	 0.7153	 0.4010
T	 0.7129	 0.3980
U	 0.6681	 0.3800
V	 0.5207	 0.3460
W	 0.7474	 0.4280
X	 0.6682	 0.3760
Y	 0.7146	 0.3890
Z	 0.7406	 0.4020



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Chain	Atom inclusion	Q-score
a	 0.6840	 0.3550
b	 0.7311	 0.4010
c	 0.6897	 0.3650
d	 0.4818	 0.3210
e	 0.4348	 0.2040
f	 0.5826	 0.3060
g	 0.6919	 0.3670
h	 0.5324	 0.3390
i	 0.7208	 0.3890
j	 0.7102	 0.3910
k	 0.5224	 0.2910
l	 0.4146	 0.2450
m	 0.4640	 0.2120
o	 0.7418	 0.3940
p	 0.6331	 0.3460
q	 0.5707	 0.3440
r	 0.6804	 0.3460
s	 0.7228	 0.3850