



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 02:51 PM UTC

PDB ID : 7PNT / pdb_00007pnt
EMDB ID : EMD-13551
Title : Assembly intermediate of mouse mitochondrial ribosome small subunit without mS37 in complex with RbfA and Tfb1m
Authors : Itoh, Y.; Khawaja, A.; Laptev, I.; Sergiev, P.; Rorbach, J.; Amunts, A.
Deposited on : 2021-09-08
Resolution : 3.19 Å(reported)
Based on initial model : 6RW4

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

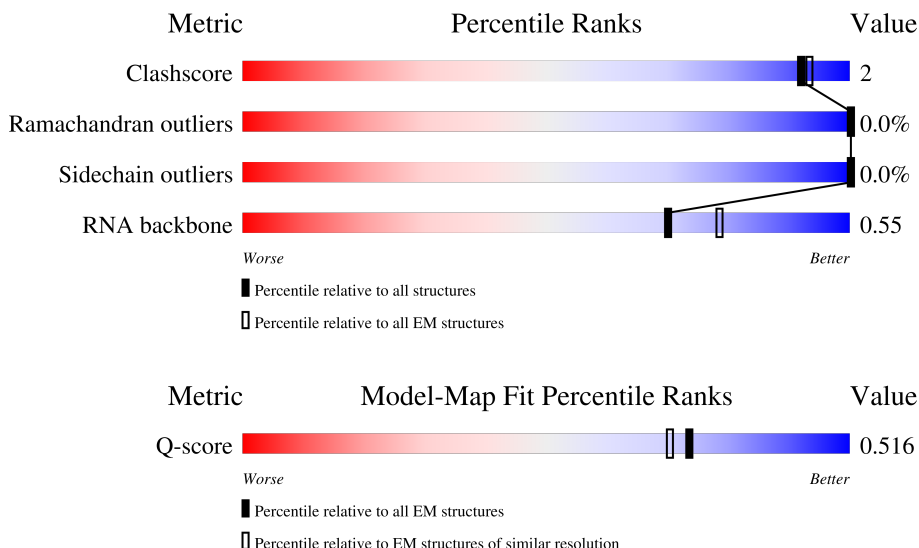
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.19 Å.

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



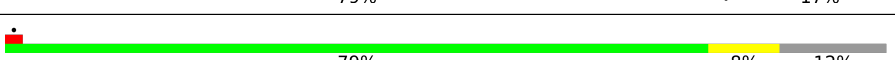
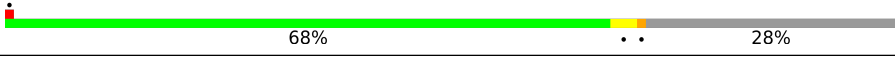
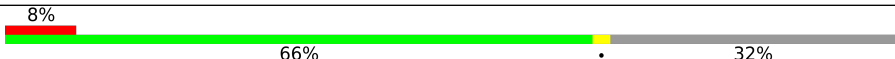
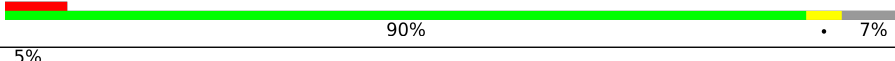
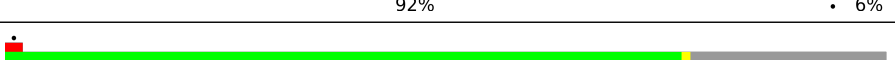
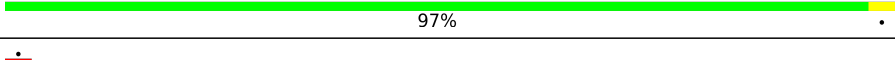
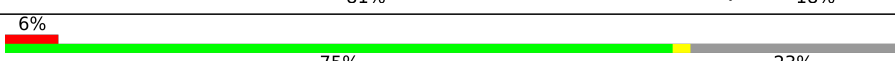
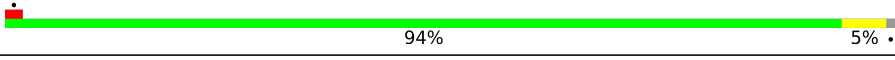
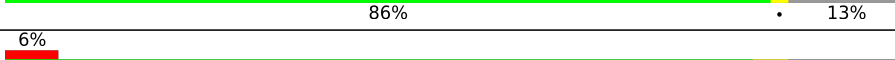
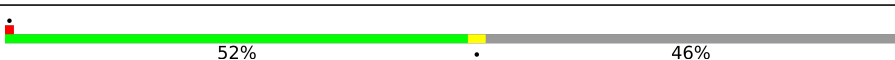
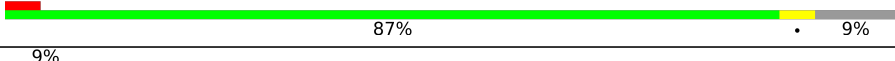
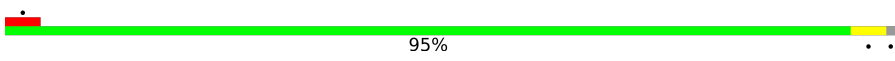
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	14455 (2.69 - 3.69)

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	A	956	
2	B	291	
3	C	167	

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Mol	Chain	Length	Quality of chain
4	D	432	
5	E	125	
6	F	242	
7	G	390	
8	H	160	
9	I	191	
10	J	139	
11	K	128	
12	L	258	
13	M	135	
14	N	120	
15	O	254	
16	P	143	
17	Q	86	
18	R	359	
19	S	177	
20	T	171	
21	U	200	
22	V	415	
23	W	186	
24	X	391	
25	Y	384	
26	Z	106	
27	0	218	
28	1	320	

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Mol	Chain	Length	Quality of chain
29	3	200	17% 24% 75%
30	4	685	29% 83% 14%
31	a	350	12% 37% 62%
32	c	345	63% 81% 6% 13%

2 Entry composition

There are 38 unique types of molecules in this entry. The entry contains 128436 atoms, of which 59523 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	A	909	Total	C	H	N	O	P	0	0
			29056	8669	9755	3459	6264	909		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	223	Total	C	H	N	O	S	0	0
			3590	1142	1799	326	315	8		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	132	Total	C	H	N	O	S	0	0
			2163	690	1091	197	180	5		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	337	Total	C	H	N	O	S	0	0
			5414	1686	2732	513	473	10		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
5	E	122	Total	C	H	N	O	S	0	0
			1986	617	1007	181	178	3		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
6	F	208	Total	C	H	N	O	S	0	0
			3472	1096	1750	316	299	11		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
7	G	325	Total	C	H	N	O	S	0	0
			5304	1689	2630	480	491	14		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
8	H	140	Total	C	H	N	O	S	0	0
			2346	742	1193	200	207	4		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
9	I	137	Total	C	H	N	O	S	0	0
			2047	629	1038	191	184	5		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	181	5F0	ASN	conflict	UNP Q9DCA2

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

Mol	Chain	Residues	Atoms						AltConf	Trace
10	J	108	Total	C	H	N	O	S	0	0
			1749	528	903	172	141	5		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
11	K	101	Total	C	H	N	O	S	0	0
			1743	534	888	175	140	6		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
12	L	176	Total	C	H	N	O	S	0	0
			3041	930	1576	274	255	6		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
13	M	126	Total	C	H	N	O	S	0	0
			2004	623	1009	194	172	6		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
14	N	113	Total	C	H	N	O	S	0	0
			1839	575	951	160	150	3		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
15	O	197	Total	C	H	N	O	S	0	0
			3146	1014	1548	289	286	9		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
16	P	97	Total	C	H	N	O	S	0	0
			1610	505	818	140	139	8		

- Molecule 17 is a protein called 28S ribosomal protein S21, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
17	Q	86	Total	C	H	N	O	S	0	0
			1482	453	750	146	126	7		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
18	R	294	Total	C	H	N	O	S	0	0
			4816	1526	2416	418	449	7		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
19	S	136	Total	C	H	N	O	S	0	0
			2257	722	1133	199	201	2		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
20	T	170	Total	C	H	N	O	S	0	0
			2801	892	1413	238	246	12		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
21	U	174	Total	C	H	N	O	S	0	0
			2908	894	1459	283	270	2		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
22	V	365	Total	C	H	N	O	S	0	0
			5970	1911	2972	506	570	11		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
23	W	100	Total	C	H	N	O	S	0	0
			1606	503	813	141	146	3		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
24	X	357	Total	C	H	N	O	S	0	0
			5762	1834	2881	515	522	10		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
25	Y	149	Total	C	H	N	O	S	0	0
			2439	809	1193	201	233	3		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
26	Z	101	Total	C	H	N	O	S	0	0
			1682	526	848	157	148	3		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
27	0	216	Total	C	H	N	O	S	0	0
			3649	1139	1838	355	313	4		

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

Mol	Chain	Residues	Atoms						AltConf	Trace
28	1	278	Total	C	H	N	O	S	0	0
			4464	1403	2241	384	424	12		

- Molecule 29 is a protein called Aurora kinase A-interacting protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
29	3	50	Total	C	H	N	O		0	0
			929	286	489	86	68			

- Molecule 30 is a protein called Pentatricopeptide repeat domain-containing protein 3, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
30	4	589	Total	C	H	N	O	S	0	0
			9558	3064	4800	800	872	22		

- Molecule 31 is a protein called Putative ribosome-binding factor A, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
31	a	134	Total	C	H	N	O	S	0	0
			2169	680	1099	190	197	3		

- Molecule 32 is a protein called Dimethyladenosine transferase 1, mitochondrial.

Mol	Chain	Residues	Atoms						AltConf	Trace
32	c	299	Total	C	H	N	O	S	0	0
			4871	1543	2478	419	421	10		

- Molecule 33 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
33	A	32	Total	Mg	0
			32	32	
33	B	1	Total	Mg	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
33	X	1	Total	Mg	0
			1	1	

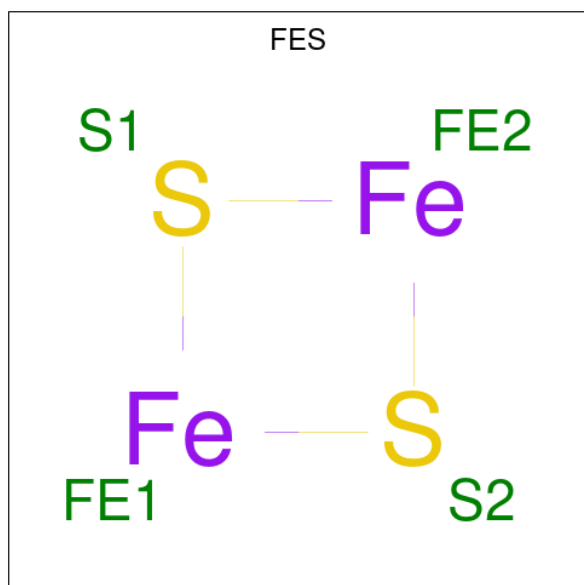
- Molecule 34 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
34	A	7	Total	K	0
			7	7	

- Molecule 35 is ZINC ION (CCD ID: ZN) (formula: Zn).

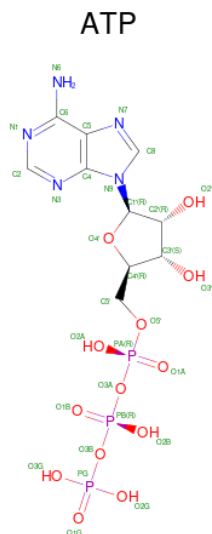
Mol	Chain	Residues	Atoms		AltConf
35	O	1	Total	Zn	0
			1	1	

- Molecule 36 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
36	P	1	Total	Fe	S	0
			4	2	2	
36	T	1	Total	Fe	S	0
			4	2	2	

- Molecule 37 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

- Molecule 38 is water.

Mol	Chain	Residues	Atoms	AltConf
38	A	268	Total O 268 268	0
38	B	7	Total O 7 7	0
38	C	23	Total O 23 23	0
38	D	14	Total O 14 14	0
38	F	5	Total O 5 5	0
38	G	9	Total O 9 9	0
38	H	11	Total O 11 11	0
38	I	3	Total O 3 3	0
38	J	4	Total O 4 4	0
38	K	15	Total O 15 15	0
38	L	1	Total O 1 1	0

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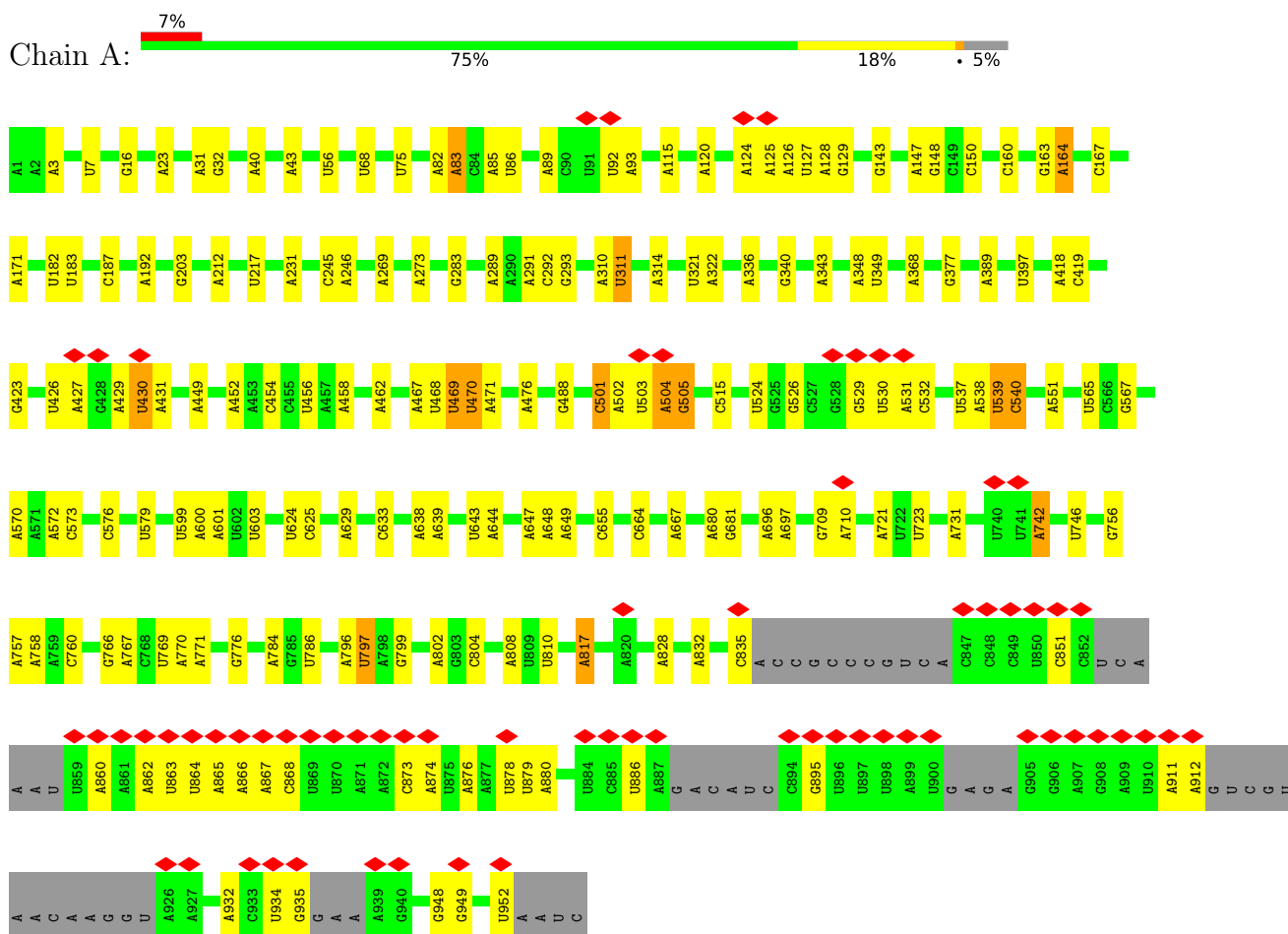
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Mol	Chain	Residues	Atoms		AltConf
38	M	17	Total 17	O 17	0
38	N	2	Total 2	O 2	0
38	O	20	Total 20	O 20	0
38	P	6	Total 6	O 6	0
38	Q	2	Total 2	O 2	0
38	R	5	Total 5	O 5	0
38	S	2	Total 2	O 2	0
38	T	6	Total 6	O 6	0
38	U	3	Total 3	O 3	0
38	X	4	Total 4	O 4	0
38	Y	3	Total 3	O 3	0
38	Z	17	Total 17	O 17	0
38	0	5	Total 5	O 5	0
38	1	11	Total 11	O 11	0
38	4	7	Total 7	O 7	0

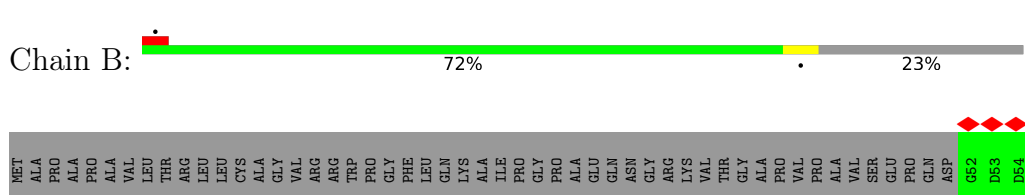
3 Residue-property plots [i](#)

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

• Molecule 1: 12S mitochondrial rRNA

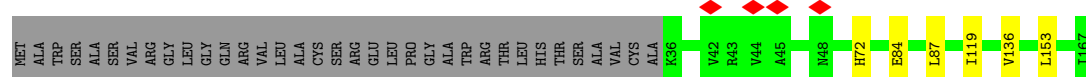
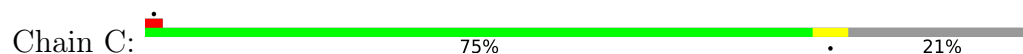


• Molecule 2: 28S ribosomal protein S2, mitochondrial

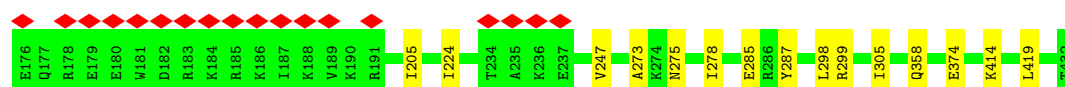
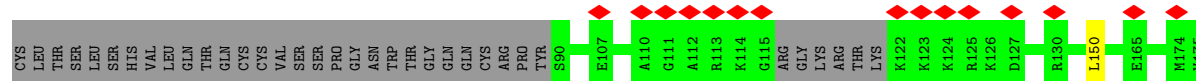
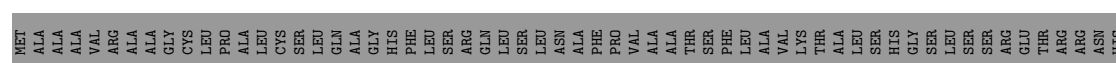




- Molecule 3: 28S ribosomal protein S24, mitochondrial



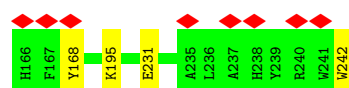
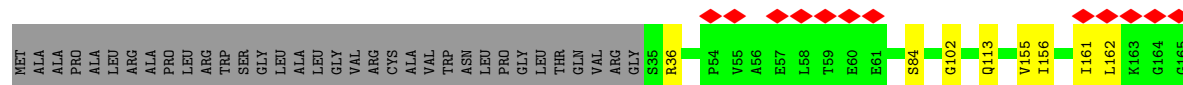
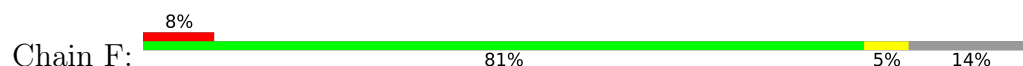
- Molecule 4: 28S ribosomal protein S5, mitochondrial



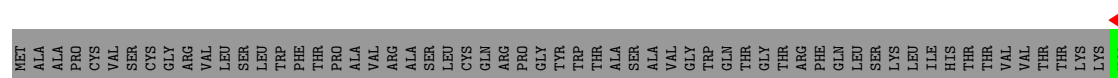
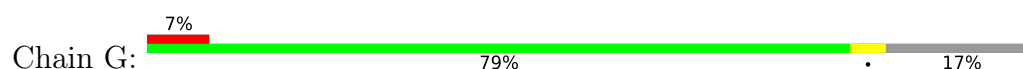
- Molecule 5: 28S ribosomal protein S6, mitochondrial

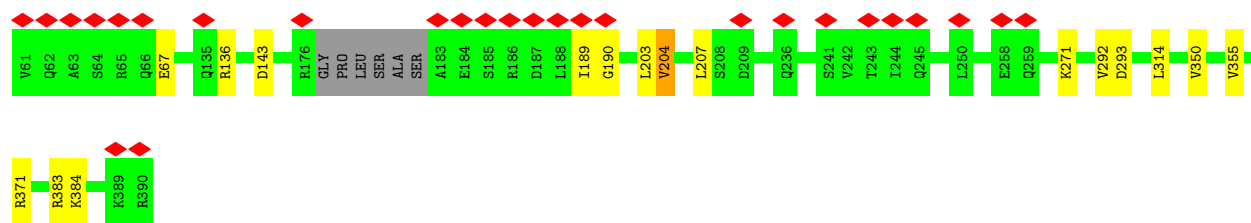


- Molecule 6: 28S ribosomal protein S7, mitochondrial

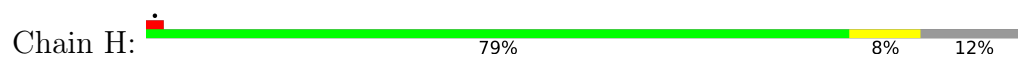


- Molecule 7: 28S ribosomal protein S9, mitochondrial

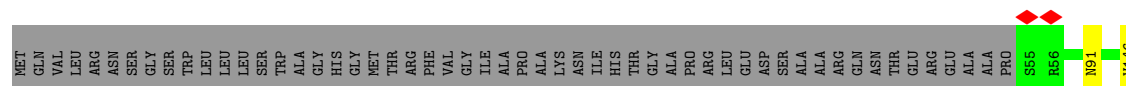




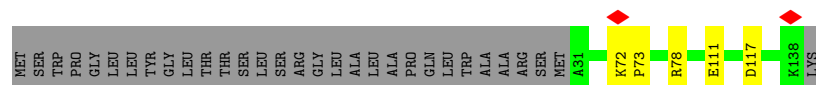
- Molecule 8: 28S ribosomal protein S10, mitochondrial



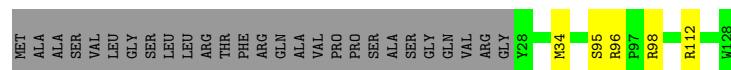
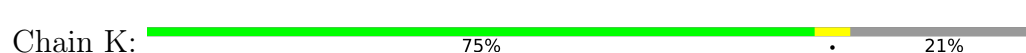
- Molecule 9: 28S ribosomal protein S11, mitochondrial



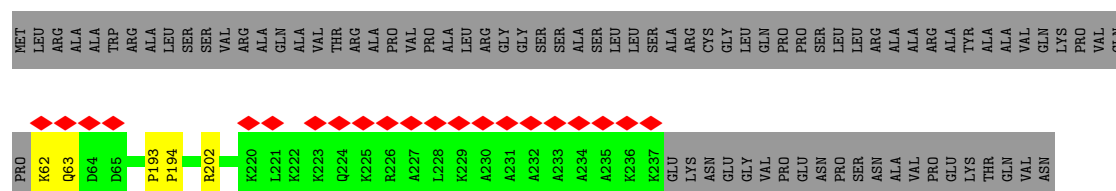
- Molecule 10: 28S ribosomal protein S12, mitochondrial



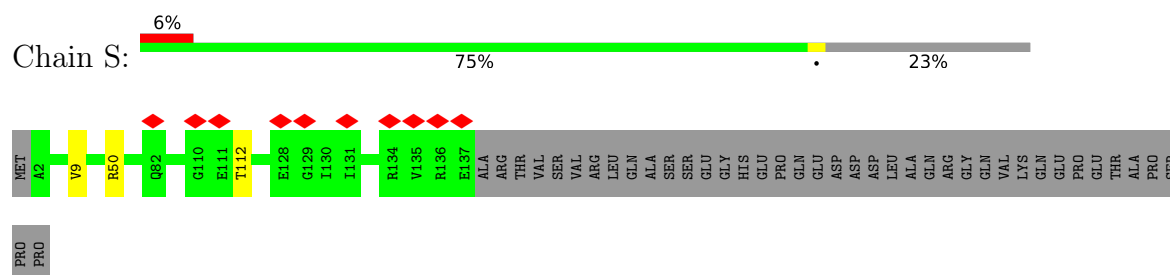
- Molecule 11: 28S ribosomal protein S14, mitochondrial



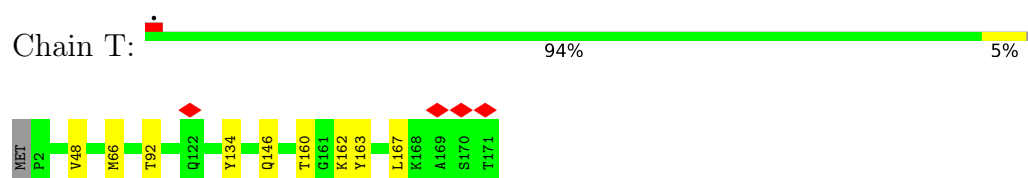
- Molecule 12: 28S ribosomal protein S15, mitochondrial



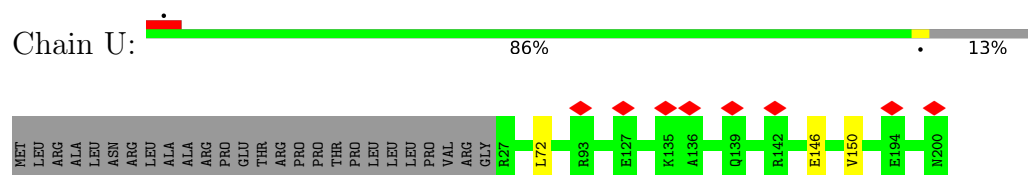
- Molecule 19: 28S ribosomal protein S23, mitochondrial



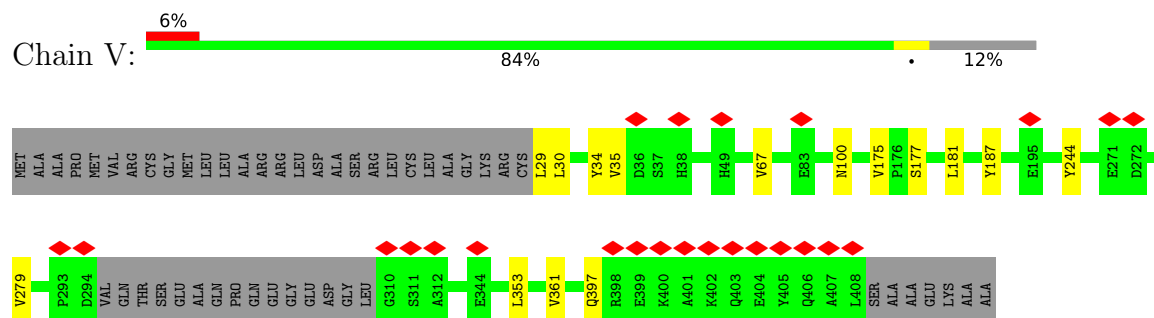
- Molecule 20: 28S ribosomal protein S25, mitochondrial



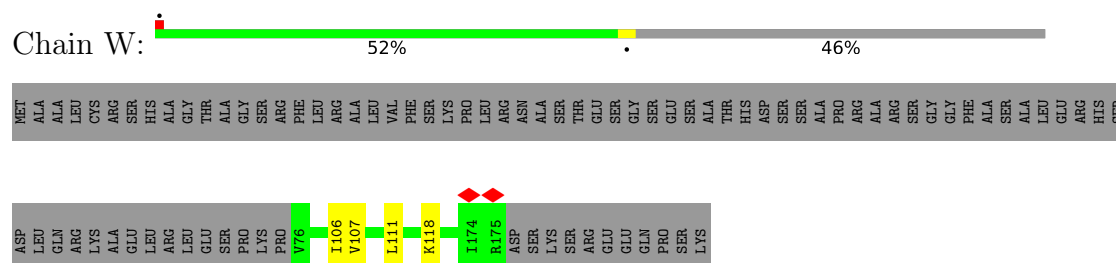
- Molecule 21: 28S ribosomal protein S26, mitochondrial



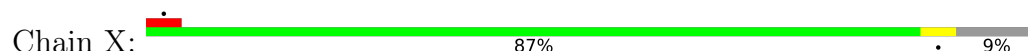
- Molecule 22: 28S ribosomal protein S27, mitochondrial



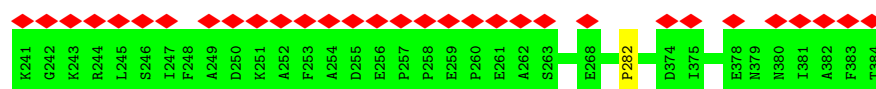
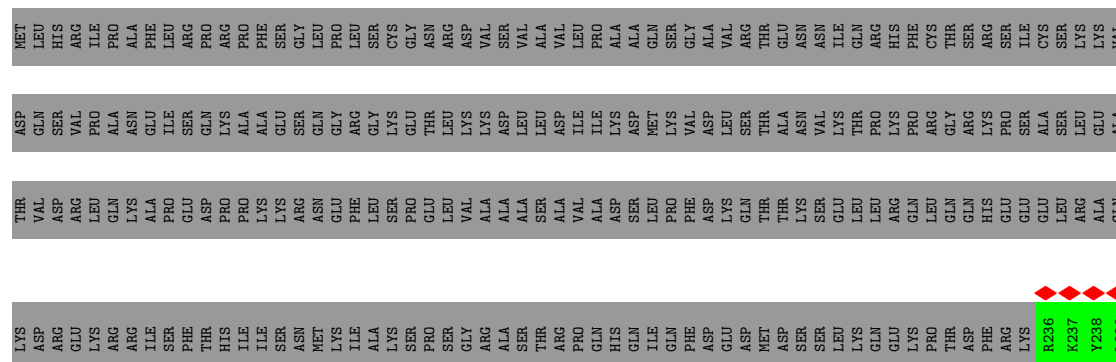
- Molecule 23: 28S ribosomal protein S28, mitochondrial



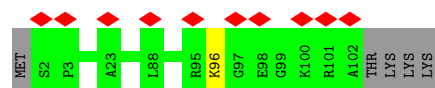
- Molecule 24: 28S ribosomal protein S29, mitochondrial



- Molecule 25: 28S ribosomal protein S31, mitochondrial



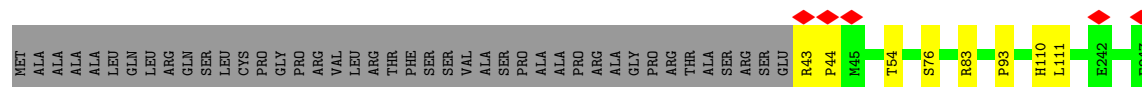
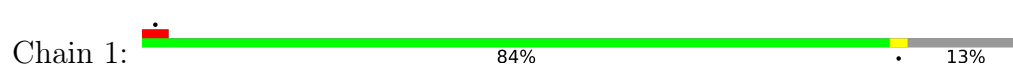
- Molecule 26: 28S ribosomal protein S33, mitochondrial

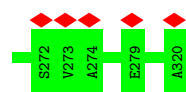


- Molecule 27: 28S ribosomal protein S34, mitochondrial



- Molecule 28: 28S ribosomal protein S35, mitochondrial





- Molecule 29: Aurora kinase A-interacting protein



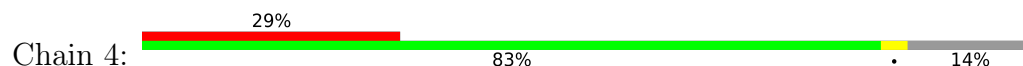
MET PHE LEU LEU LEU THR SER ARG ARG ALA ALA THR THR VAL VAL VAL PRO TRP ALA GLY PHE SER ARG ARG CYS PRO GLY SER ARG GLY VAL ILE ILE SER TYR ALA PHE THR LEU ARG PRO LEU TTR SER GLN LEU PRO ALA SER PRO ARG GLN ARG ALA ALA SER LEU PRO GLN GLY THR ARG GLN SER GLU

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

T151 R152 F153 L154 R155 R156 K157 V158 R159 E160 G161 R162 L163 K164 K165 I168 K172 K175 L179 K180 A181 G182 L183 K184 E185 A186

P187 E188 N189 W190 Q191 T192 P193 K194 I195 Y196 L197 K198 N199 K200

- Molecule 30: Pentatricopeptide repeat domain-containing protein 3, mitochondrial



MET ALA ALA ALA VAL ALA VAL SER PHE ARG SER GLY VAL VAL CYS GLU PRO LYS VAL CYS CYS ARG PHE TYR ALA GLY THR GLU SER LEU PRO LYS VAL ILE E56 E57 I58 V59

V89 L109 E147 P148 Q149 I150 E151 D152 V153 S154 E155 A156 E159 E160 R161 R165 K166 V167 R168 D172 M173 F174 D175 Q176 L177 L178 L179 Q179 A180 G181 T182 T183 V184 S185 L186 E187 T188 T189 L193 Y198 Y199 G200 D201 Q202 P204 P205 A206 D207 Y208 P209 PHE GLN

THR GLU HIS LEU ASN LEU GLU ALA ALA GLU GLU ALA ALA GLU ASN ASN THR SER LYS MET GLU SER G235 P236 W237 K238 A239 Q240 N241 N242 A243 E244 R245 T246 F247 A248 L249 W250 P251 E252 K253 H256 R269 A270 Y271 A272 Q273 A274 L275 N276 V277 Y278 T279 E280 L281 L282 N283 N284 E285

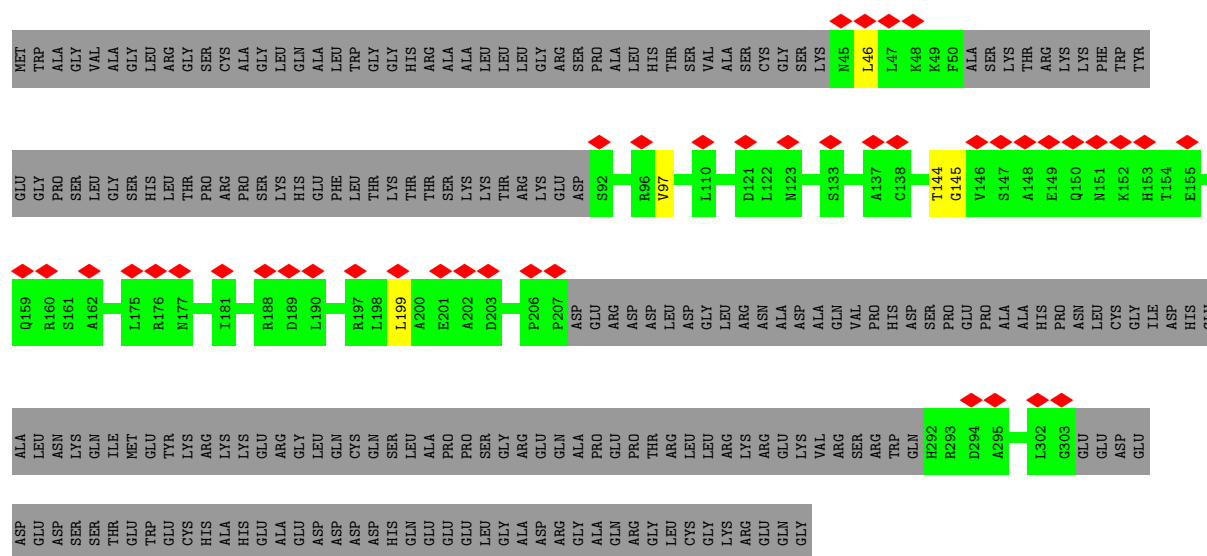
L286 S287 A288 D289 N294 I297 K300 T301 F302 L303 A304 N305 E306 K307 F308 E309 K311 W312 N313 I315 L316 D317 L318 L319 K320 H321 K322 V327 K328 P329 M330 L331 K343 C344 Y345 S346 L347 G348 R349 I350 P351 A352 V353 L356 R357 E358 K359 K360 H361 I362 G363 I364

E365 L368 Y371 F378 Y379 P380 R381 D382 L383 S384 A385 I386 K387 M388 P389 S390 L391 I392 E399 L400 Q408 D409 L410 D411 L447 H450 E469 P567 A568 F569 D570 W571 P572 P575 G587 R588 S589 Q590 E591 A592 W593 K594 E597 K600 K601 I602 E603

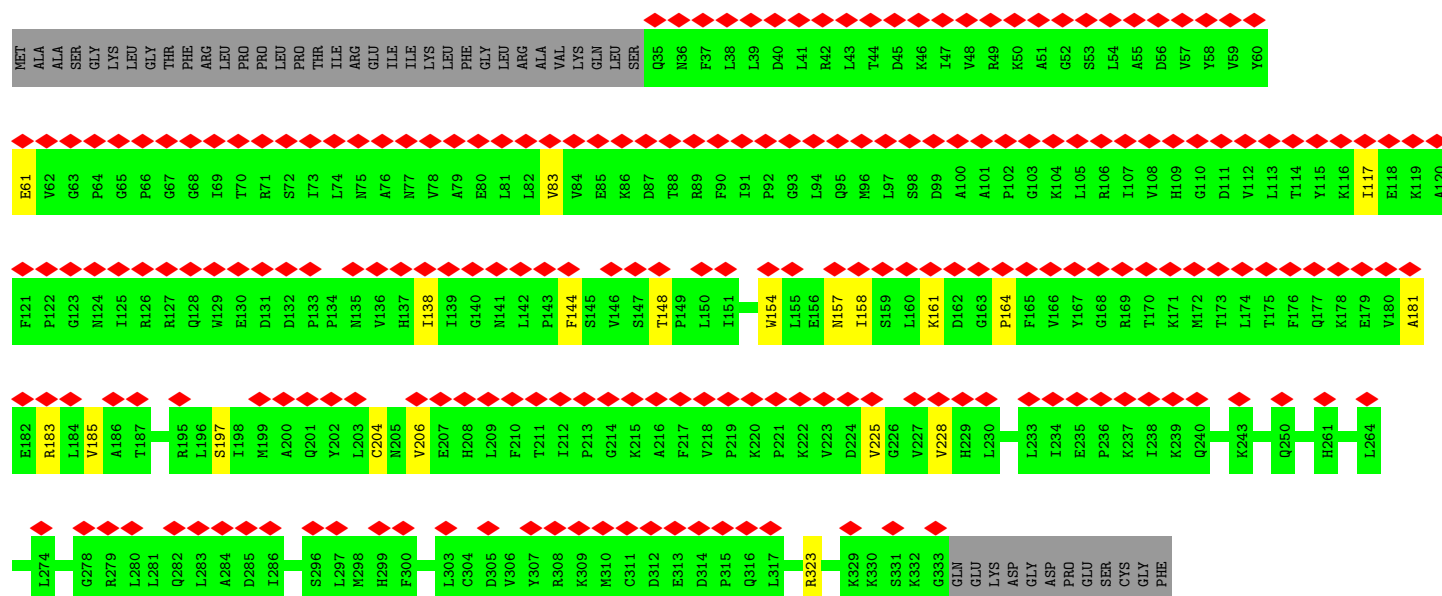
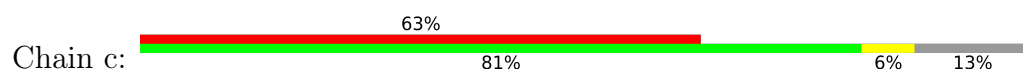
K604 R607 M615 D616 T617 A618 K619 A620 S621 G622 S623 T624 A625 L626 A627 T628 E629 V630 V631 K632 L633 A634 S635 A636 F637 S638 L639 P640 I641 G642 E643 S644 L645 A646 Q647 R648 V649 V650 M651 D652 F653 T654 V655 D656 P657 E658 Q659 K660 E661 A662 L663 G664 N665 L666 T667 E668 I669 ASN

SER SER ASP GLY SER SER ASP ASP ASP ASP LYS

- Molecule 31: Putative ribosome-binding factor A, mitochondrial



- Molecule 32: Dimethyladenosine transferase 1, mitochondrial



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	52361	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 ($e^-/\text{\AA}^2$)	31	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	3600	Depositor
Magnification	165000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	1.622	Depositor
Minimum map value	-0.805	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	398.40192, 398.40192, 398.40192	wwPDB
Map dimensions	576, 576, 576	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.69167, 0.69167, 0.69167	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, ZN, ATP, MG, K, 5F0, AYA

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	A	0.10	0/21600	0.19	0/33606
2	B	0.13	0/1832	0.30	0/2479
3	C	0.14	0/1100	0.29	0/1485
4	D	0.12	0/2737	0.28	0/3671
5	E	0.12	0/997	0.31	0/1347
6	F	0.11	0/1763	0.28	0/2368
7	G	0.11	0/2734	0.29	0/3669
8	H	0.14	0/1179	0.28	0/1597
9	I	0.13	0/1018	0.33	0/1374
10	J	0.14	0/862	0.34	0/1155
11	K	0.15	0/871	0.34	0/1167
12	L	0.11	0/1485	0.26	0/1980
13	M	0.15	0/1017	0.32	0/1366
14	N	0.13	0/907	0.31	0/1228
15	O	0.12	0/1653	0.29	0/2254
16	P	0.13	0/809	0.29	0/1085
17	Q	0.12	0/735	0.33	0/980
18	R	0.12	0/2449	0.27	0/3311
19	S	0.13	0/1148	0.27	0/1541
20	T	0.12	0/1420	0.28	0/1903
21	U	0.13	0/1470	0.29	0/1976
22	V	0.09	0/3059	0.24	0/4135
23	W	0.11	0/805	0.29	0/1084
24	X	0.10	0/2952	0.27	0/3995
25	Y	0.11	0/1283	0.25	0/1730
26	Z	0.13	0/851	0.28	0/1133
27	0	0.13	0/1856	0.33	0/2511
28	1	0.11	0/2271	0.27	0/3078
29	3	0.09	0/448	0.29	0/591
30	4	0.10	0/4868	0.25	0/6597
31	a	0.08	0/1087	0.24	0/1470
32	c	0.09	0/2447	0.25	0/3309

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.11	0/71713	0.25	0/101175

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
9	I	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
9	I	181	5F0	Mainchain

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	A	19301	9755	9761	45	0
2	B	1791	1799	1797	9	0
3	C	1072	1091	1087	5	0
4	D	2682	2732	2727	12	0
5	E	979	1007	1007	6	0
6	F	1722	1750	1748	10	0
7	G	2674	2630	2626	11	0
8	H	1153	1193	1190	9	0
9	I	1009	1038	1029	4	0
10	J	846	903	901	3	0
11	K	855	888	887	3	0
12	L	1465	1576	1574	3	0
13	M	995	1009	1006	4	0
14	N	888	951	947	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	O	1598	1548	1547	3	0
16	P	792	818	817	6	0
17	Q	732	750	750	3	0
18	R	2400	2416	2415	3	0
19	S	1124	1133	1132	3	0
20	T	1388	1413	1413	7	0
21	U	1449	1459	1456	2	0
22	V	2998	2972	2967	11	0
23	W	793	813	811	3	0
24	X	2881	2881	2879	10	0
25	Y	1246	1193	1191	1	0
26	Z	834	848	847	1	0
27	0	1811	1838	1834	5	0
28	1	2223	2241	2238	4	0
29	3	440	489	488	1	0
30	4	4758	4800	4793	13	0
31	a	1070	1099	1096	4	0
32	c	2393	2478	2478	13	0
33	A	32	0	0	0	0
33	B	1	0	0	0	0
33	X	1	0	0	0	0
34	A	7	0	0	0	0
35	O	1	0	0	0	0
36	P	4	0	0	0	0
36	T	4	0	0	0	0
37	X	31	12	12	0	0
38	0	5	0	0	1	0
38	1	11	0	0	0	0
38	4	7	0	0	0	0
38	A	268	0	0	4	0
38	B	7	0	0	0	0
38	C	23	0	0	1	0
38	D	14	0	0	0	0
38	F	5	0	0	2	0
38	G	9	0	0	1	0
38	H	11	0	0	0	0
38	I	3	0	0	0	0
38	J	4	0	0	1	0
38	K	15	0	0	0	0
38	L	1	0	0	0	0
38	M	17	0	0	0	0
38	N	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
38	O	20	0	0	2	0
38	P	6	0	0	0	0
38	Q	2	0	0	0	0
38	R	5	0	0	0	0
38	S	2	0	0	0	0
38	T	6	0	0	0	0
38	U	3	0	0	0	0
38	X	4	0	0	0	0
38	Y	3	0	0	0	0
38	Z	17	0	0	0	0
All	All	68913	59523	59451	181	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.

All (181) 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:A:567:G:N2	1:A:576:C:O2	2.20	0.74
10:J:111:GLU:OE1	38:J:201:HOH:O	2.05	0.74
1:A:217:U:OP1	15:O:146:LYS:NZ	2.22	0.73
1:A:192:A:N7	38:A:1105:HOH:O	2.21	0.72
6:F:84:SER:OG	24:X:372:GLU:OE2	2.07	0.72
32:c:117:ILE:HD11	32:c:164:PRO:HB3	1.72	0.71
3:C:72:HIS:O	38:C:201:HOH:O	2.08	0.71
1:A:336:A:OP1	9:I:91:ASN:ND2	2.24	0.71
7:G:271:LYS:O	38:G:401:HOH:O	2.12	0.67
7:G:383:ARG:NH2	7:G:384:LYS:O	2.28	0.67
32:c:158:ILE:HD13	32:c:204:CYS:SG	2.36	0.66
1:A:935:G:N7	32:c:183:ARG:NE	2.42	0.66
8:H:51:GLU:OE1	8:H:100:ARG:NH1	2.27	0.66
1:A:449:A:OP1	1:A:505:G:N2	2.28	0.66
13:M:55:ASP:OD2	20:T:146:GLN:NE2	2.29	0.66
7:G:203:LEU:HD12	7:G:207:LEU:HD11	1.79	0.63
32:c:61:GLU:HB3	32:c:83:VAL:HG12	1.83	0.61
24:X:176:GLU:N	24:X:176:GLU:OE2	2.33	0.61
20:T:92:THR:HG22	20:T:92:THR:O	2.01	0.60
6:F:161:ILE:O	6:F:168:TYR:N	2.35	0.60
1:A:647:A:OP1	2:B:198:ASN:ND2	2.34	0.60
24:X:116:ARG:NH2	24:X:330:LEU:O	2.35	0.59
11:K:96:ARG:NH2	11:K:98:ARG:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:804:C:OP2	38:A:1102:HOH:O	2.16	0.59
22:V:279:VAL:HG23	22:V:353:LEU:HD11	1.85	0.59
32:c:144:PHE:HZ	32:c:225:VAL:HG22	1.68	0.59
14:N:3:ILE:HG22	14:N:3:ILE:O	2.02	0.59
1:A:723:U:OP1	6:F:195:LYS:NZ	2.35	0.58
15:O:136:PRO:O	38:O:402:HOH:O	2.17	0.58
5:E:94:VAL:HG11	16:P:118:MET:HE1	1.85	0.58
1:A:501:C:OP1	12:L:202:ARG:NH2	2.35	0.58
1:A:817:A:O2'	6:F:242:TRP:OXT	2.22	0.57
9:I:148:VAL:HG13	9:I:148:VAL:O	2.03	0.57
1:A:504:A:O2'	29:3:152:ARG:NH1	2.37	0.57
22:V:29:LEU:HD12	22:V:181:LEU:HD23	1.85	0.57
4:D:287:TYR:OH	4:D:374:GLU:OE2	2.22	0.56
22:V:279:VAL:CG2	22:V:353:LEU:HD11	2.36	0.55
1:A:551:A:HO2'	1:A:771:A:HO2'	1.53	0.55
1:A:746:U:O2'	1:A:802:A:N6	2.39	0.55
22:V:67:VAL:O	22:V:100:ASN:ND2	2.39	0.55
22:V:35:VAL:HG12	22:V:35:VAL:O	2.07	0.55
1:A:696:A:OP2	26:Z:96:LYS:NZ	2.40	0.54
1:A:633:C:O2'	1:A:649:A:N1	2.39	0.54
32:c:157:ASN:O	32:c:161:LYS:N	2.40	0.54
1:A:469:U:OP2	19:S:50:ARG:NH2	2.41	0.53
1:A:599:U:O2'	1:A:600:A:N7	2.39	0.53
6:F:36:ARG:NH2	38:F:302:HOH:O	2.40	0.53
28:1:83:ARG:NH1	28:1:93:PRO:O	2.39	0.53
5:E:117:LEU:HD13	31:a:46:LEU:HD22	1.91	0.53
18:R:161:SER:O	18:R:169:ARG:NH1	2.40	0.53
1:A:31:A:O3'	38:A:1101:HOH:O	2.18	0.52
5:E:94:VAL:CG1	16:P:118:MET:HE1	2.40	0.52
16:P:68:LEU:HD21	16:P:80:LEU:HD11	1.90	0.52
10:J:78:ARG:NE	10:J:117:ASP:OD2	2.43	0.52
20:T:160:THR:HG22	20:T:162:LYS:H	1.75	0.51
1:A:43:A:N7	1:A:68:U:O2'	2.42	0.51
3:C:119:ILE:HB	3:C:153:LEU:HD12	1.92	0.51
6:F:113:GLN:OE1	38:F:301:HOH:O	2.19	0.51
2:B:86:VAL:HG12	2:B:86:VAL:O	2.11	0.51
32:c:185:VAL:HG12	32:c:185:VAL:O	2.11	0.51
1:A:579:U:O2'	1:A:796:A:O4'	2.28	0.51
4:D:285:GLU:O	4:D:358:GLN:NE2	2.44	0.51
30:4:386:ILE:HG22	30:4:386:ILE:O	2.11	0.51
7:G:292:VAL:HG12	7:G:293:ASP:N	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:150:LEU:HD21	30:4:109:LEU:CB	2.41	0.49
24:X:41:VAL:HG12	24:X:41:VAL:O	2.10	0.49
1:A:430:U:N3	32:c:148:THR:OG1	2.43	0.49
8:H:80:LEU:HD11	8:H:87:LYS:HA	1.94	0.49
27:0:103:ASP:N	27:0:111:HIS:O	2.43	0.49
21:U:146:GLU:O	21:U:150:VAL:HG23	2.12	0.48
2:B:272:SER:HB3	2:B:273:PRO:HD3	1.94	0.48
2:B:236:ASN:OD1	23:W:118:LYS:NZ	2.47	0.48
11:K:34:MET:HE3	11:K:95:SER:HB2	1.96	0.48
22:V:187:TYR:CE1	22:V:353:LEU:HD13	2.49	0.48
13:M:28:ASN:ND2	20:T:160:THR:HG23	2.29	0.48
24:X:144:LEU:CD2	24:X:240:LEU:HD22	2.44	0.48
25:Y:282:PRO:HB2	30:4:89:VAL:CG1	2.44	0.48
4:D:275:ASN:O	4:D:278:ILE:HG22	2.14	0.48
28:1:54:THR:HG23	28:1:76:SER:O	2.14	0.47
1:A:470:U:O2'	4:D:299:ARG:NH1	2.48	0.47
30:4:59:VAL:O	30:4:59:VAL:HG23	2.14	0.47
5:E:79:LEU:HG	5:E:93:ILE:HD12	1.96	0.47
8:H:39:VAL:HG22	8:H:100:ARG:O	2.15	0.47
32:c:183:ARG:O	32:c:197:SER:OG	2.22	0.47
1:A:147:A:OP2	20:T:134:TYR:OH	2.31	0.47
1:A:127:U:O4	1:A:128:A:N6	2.48	0.46
1:A:321:U:O2'	1:A:322:A:N7	2.40	0.46
27:0:61:GLU:O	38:0:301:HOH:O	2.21	0.46
1:A:163:G:H2'	1:A:164:A:H5''	1.98	0.46
1:A:539:U:O2	1:A:540:C:N4	2.49	0.46
20:T:48:VAL:HG13	20:T:66:MET:SD	2.56	0.46
4:D:414:LYS:HG3	4:D:419:LEU:HD22	1.97	0.46
1:A:504:A:N6	1:A:932:A:OP1	2.44	0.46
1:A:797:U:O2'	1:A:799:G:N7	2.48	0.46
1:A:419:C:O2	32:c:323:ARG:NH2	2.36	0.46
1:A:601:A:OP2	38:A:1104:HOH:O	2.21	0.46
10:J:72:LYS:CG	10:J:73:PRO:HA	2.46	0.46
8:H:29:ASP:OD1	8:H:30:THR:N	2.48	0.46
1:A:808:A:OP2	7:G:371:ARG:NH2	2.48	0.46
8:H:81:LYS:O	11:K:112:ARG:NH2	2.49	0.46
1:A:340:G:C5	31:a:97:VAL:HG22	2.51	0.45
5:E:54:HIS:NE2	5:E:85:ASP:O	2.49	0.45
24:X:92:LEU:HD21	24:X:133:HIS:HB2	1.98	0.45
1:A:160:C:OP1	27:0:19:ARG:NH2	2.50	0.45
8:H:104:LEU:HD12	8:H:104:LEU:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:88:VAL:HG13	14:N:88:VAL:O	2.16	0.45
16:P:139:ILE:HG23	16:P:139:ILE:O	2.17	0.45
4:D:298:LEU:C	4:D:298:LEU:HD12	2.41	0.45
6:F:155:VAL:HG23	6:F:156:ILE:N	2.31	0.45
30:4:371:TYR:CE2	30:4:400:LEU:HD21	2.52	0.45
31:a:144:THR:HG22	31:a:145:GLY:N	2.32	0.45
1:A:766:G:OP1	24:X:272:ARG:NH2	2.48	0.45
32:c:181:ALA:HB1	32:c:228:VAL:CG2	2.47	0.45
19:S:9:VAL:O	19:S:9:VAL:HG13	2.17	0.45
7:G:314:LEU:HD13	7:G:350:VAL:HG11	1.98	0.44
32:c:138:ILE:HG21	32:c:154:TRP:CH2	2.52	0.44
3:C:87:LEU:HD13	3:C:87:LEU:O	2.18	0.44
12:L:62:LYS:HG2	12:L:63:GLN:H	1.81	0.44
18:R:275:VAL:HG11	18:R:306:LEU:HD12	2.00	0.44
22:V:29:LEU:HD13	22:V:361:VAL:HG11	1.98	0.44
30:4:368:LEU:HD12	30:4:447:LEU:HD13	1.99	0.44
2:B:75:LEU:HD21	2:B:256:ARG:CZ	2.48	0.44
15:O:91:THR:OG1	38:O:401:HOH:O	2.16	0.44
17:Q:42:ARG:O	17:Q:42:ARG:HG3	2.18	0.44
22:V:175:VAL:HG12	22:V:177:SER:H	1.83	0.44
24:X:275:ILE:HG22	24:X:276:ALA:N	2.32	0.44
24:X:265:THR:OG1	24:X:275:ILE:O	2.36	0.44
30:4:350:ILE:HB	30:4:351:PRO:HD3	2.00	0.44
1:A:82:A:H2'	1:A:83:A:O4'	2.17	0.43
1:A:171:A:O2'	1:A:183:U:O2'	2.28	0.43
17:Q:83:PRO:HA	23:W:107:VAL:HG21	1.99	0.43
1:A:876:A:OP2	22:V:397:GLN:NE2	2.51	0.43
2:B:185:ARG:NH2	7:G:143:ASP:O	2.48	0.43
27:O:37:ASP:O	27:O:41:LEU:N	2.52	0.43
6:F:231:GLU:HA	31:a:199:LEU:HD11	1.99	0.43
22:V:30:LEU:HD13	22:V:34:TYR:CD1	2.54	0.43
8:H:24:THR:O	8:H:24:THR:HG23	2.18	0.43
13:M:104:ILE:HG23	18:R:146:ILE:HG12	2.00	0.43
4:D:205:ILE:HG21	4:D:224:ILE:HD11	2.00	0.43
16:P:114:ARG:HG3	16:P:118:MET:HE3	2.01	0.43
24:X:146:LEU:HD21	24:X:237:LEU:HD22	2.01	0.43
16:P:139:ILE:HD11	17:Q:32:THR:CG2	2.49	0.42
6:F:161:ILE:HG22	6:F:162:LEU:N	2.33	0.42
12:L:193:PRO:HA	12:L:194:PRO:HD3	1.96	0.42
1:A:721:A:N6	1:A:742:A:N3	2.68	0.42
1:A:310:A:OP1	1:A:311:U:N3	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:189:ILE:HG22	7:G:190:GLY:N	2.35	0.42
28:1:110:HIS:CD2	28:1:111:LEU:HG	2.55	0.42
2:B:143:SER:O	2:B:165:THR:HA	2.20	0.42
3:C:136:VAL:HG22	3:C:153:LEU:HD23	2.02	0.42
7:G:350:VAL:HG23	7:G:355:VAL:HG23	2.01	0.42
19:S:112:THR:HG22	19:S:112:THR:O	2.19	0.42
9:I:173:THR:HG22	9:I:174:ASP:N	2.35	0.42
1:A:810:U:O2'	6:F:102:GLY:O	2.38	0.42
30:4:297:ILE:HG23	30:4:315:ILE:HG23	2.02	0.42
2:B:217:VAL:HG22	2:B:231:TYR:HB2	2.02	0.41
5:E:47:LEU:HD13	5:E:51:ILE:HD12	2.01	0.41
23:W:106:ILE:HG23	23:W:111:LEU:HD23	2.01	0.41
27:O:71:LEU:HD11	27:O:141:LEU:HD22	2.01	0.41
3:C:84:GLU:OE1	3:C:84:GLU:N	2.54	0.41
4:D:150:LEU:HD21	30:4:109:LEU:HB3	2.03	0.41
20:T:163:TYR:CE2	20:T:167:LEU:HD11	2.55	0.41
1:A:89:A:O3'	21:U:72:LEU:HD13	2.21	0.41
1:A:343:A:N3	1:A:423:G:O2'	2.37	0.41
22:V:244:TYR:CD1	22:V:244:TYR:C	2.99	0.41
32:c:185:VAL:HG22	32:c:206:VAL:HG12	2.03	0.41
1:A:756:G:O2'	1:A:758:A:N7	2.48	0.41
4:D:150:LEU:HD21	30:4:109:LEU:HB2	2.02	0.41
7:G:67:GLU:O	7:G:136:ARG:NH2	2.54	0.41
8:H:104:LEU:HD21	8:H:115:TYR:CE2	2.56	0.41
30:4:167:VAL:HG23	30:4:168:ARG:N	2.36	0.41
30:4:572:PRO:HB2	30:4:575:PRO:HD2	2.02	0.41
13:M:17:LEU:HD23	13:M:81:ALA:HB2	2.02	0.41
28:1:43:ARG:N	28:1:44:PRO:CD	2.84	0.41
30:4:378:PHE:CD2	30:4:392:ILE:HD12	2.56	0.41
2:B:191:ILE:HA	2:B:217:VAL:O	2.21	0.41
9:I:146:VAL:HG23	9:I:169:VAL:HG13	2.04	0.40
4:D:298:LEU:HD11	4:D:305:ILE:HD13	2.04	0.40
8:H:104:LEU:HB3	8:H:107:LEU:HD11	2.03	0.40
7:G:204:VAL:HG12	7:G:204:VAL:O	2.22	0.40
4:D:247:VAL:HG22	4:D:273:ALA:HB1	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	221/291 (76%)	219 (99%)	2 (1%)	0	100	100
3	C	130/167 (78%)	124 (95%)	6 (5%)	0	100	100
4	D	333/432 (77%)	328 (98%)	5 (2%)	0	100	100
5	E	120/125 (96%)	118 (98%)	2 (2%)	0	100	100
6	F	206/242 (85%)	200 (97%)	6 (3%)	0	100	100
7	G	321/390 (82%)	315 (98%)	5 (2%)	1 (0%)	36	68
8	H	138/160 (86%)	135 (98%)	2 (1%)	1 (1%)	18	52
9	I	134/191 (70%)	130 (97%)	4 (3%)	0	100	100
10	J	106/139 (76%)	105 (99%)	1 (1%)	0	100	100
11	K	99/128 (77%)	98 (99%)	1 (1%)	0	100	100
12	L	174/258 (67%)	173 (99%)	1 (1%)	0	100	100
13	M	124/135 (92%)	124 (100%)	0	0	100	100
14	N	111/120 (92%)	108 (97%)	3 (3%)	0	100	100
15	O	195/254 (77%)	192 (98%)	3 (2%)	0	100	100
16	P	95/143 (66%)	94 (99%)	1 (1%)	0	100	100
17	Q	84/86 (98%)	83 (99%)	1 (1%)	0	100	100
18	R	292/359 (81%)	284 (97%)	8 (3%)	0	100	100
19	S	134/177 (76%)	132 (98%)	2 (2%)	0	100	100
20	T	168/171 (98%)	167 (99%)	1 (1%)	0	100	100
21	U	172/200 (86%)	170 (99%)	2 (1%)	0	100	100
22	V	361/415 (87%)	357 (99%)	4 (1%)	0	100	100
23	W	98/186 (53%)	96 (98%)	2 (2%)	0	100	100
24	X	355/391 (91%)	350 (99%)	5 (1%)	0	100	100
25	Y	147/384 (38%)	145 (99%)	2 (1%)	0	100	100
26	Z	99/106 (93%)	98 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
27	0	214/218 (98%)	211 (99%)	3 (1%)	0	100	100
28	1	276/320 (86%)	274 (99%)	2 (1%)	0	100	100
29	3	48/200 (24%)	48 (100%)	0	0	100	100
30	4	585/685 (85%)	579 (99%)	6 (1%)	0	100	100
31	a	128/350 (37%)	126 (98%)	2 (2%)	0	100	100
32	c	297/345 (86%)	292 (98%)	5 (2%)	0	100	100
All	All	5965/7768 (77%)	5875 (98%)	88 (2%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	H	85	ILE
7	G	204	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	195/247 (79%)	195 (100%)	0	100	100
3	C	113/139 (81%)	113 (100%)	0	100	100
4	D	277/354 (78%)	277 (100%)	0	100	100
5	E	107/110 (97%)	107 (100%)	0	100	100
6	F	180/204 (88%)	180 (100%)	0	100	100
7	G	282/335 (84%)	282 (100%)	0	100	100
8	H	130/150 (87%)	130 (100%)	0	100	100
9	I	102/143 (71%)	102 (100%)	0	100	100
10	J	93/117 (80%)	93 (100%)	0	100	100
11	K	90/110 (82%)	90 (100%)	0	100	100
12	L	161/224 (72%)	161 (100%)	0	100	100
13	M	103/112 (92%)	103 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	97/104 (93%)	97 (100%)	0	100	100
15	O	176/225 (78%)	176 (100%)	0	100	100
16	P	89/125 (71%)	89 (100%)	0	100	100
17	Q	77/77 (100%)	77 (100%)	0	100	100
18	R	261/313 (83%)	261 (100%)	0	100	100
19	S	117/152 (77%)	117 (100%)	0	100	100
20	T	153/154 (99%)	153 (100%)	0	100	100
21	U	149/171 (87%)	149 (100%)	0	100	100
22	V	326/362 (90%)	326 (100%)	0	100	100
23	W	87/156 (56%)	87 (100%)	0	100	100
24	X	314/346 (91%)	314 (100%)	0	100	100
25	Y	133/341 (39%)	133 (100%)	0	100	100
26	Z	88/93 (95%)	88 (100%)	0	100	100
27	0	190/191 (100%)	189 (100%)	1 (0%)	81	85
28	1	248/279 (89%)	248 (100%)	0	100	100
29	3	46/176 (26%)	46 (100%)	0	100	100
30	4	519/599 (87%)	519 (100%)	0	100	100
31	a	122/299 (41%)	122 (100%)	0	100	100
32	c	266/304 (88%)	266 (100%)	0	100	100
All	All	5291/6712 (79%)	5290 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
27	0	48	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	164	HIS
2	B	198	ASN
4	D	358	GLN
4	D	411	GLN
6	F	113	GLN

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Mol	Chain	Res	Type
6	F	215	ASN
7	G	288	HIS
7	G	378	GLN
9	I	84	HIS
10	J	105	HIS
12	L	119	ASN
12	L	196	HIS
13	M	75	HIS
14	N	9	HIS
15	O	96	ASN
18	R	109	GLN
18	R	304	HIS
19	S	91	ASN
20	T	51	ASN
22	V	141	ASN
22	V	145	ASN
22	V	325	GLN
22	V	397	GLN
23	W	105	HIS
24	X	136	HIS
24	X	141	HIS
24	X	228	ASN
24	X	319	HIS
25	Y	370	ASN
25	Y	379	ASN
26	Z	54	ASN
27	0	137	HIS
27	0	145	HIS
28	1	143	HIS
28	1	182	HIS
30	4	372	HIS
30	4	376	HIS
30	4	521	HIS
30	4	577	GLN
31	a	114	GLN
31	a	173	GLN
32	c	75	ASN
32	c	208	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	902/956 (94%)	138 (15%)	0

All (138) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	3	A
1	A	7	U
1	A	16	G
1	A	23	A
1	A	32	G
1	A	40	A
1	A	56	U
1	A	75	U
1	A	83	A
1	A	85	A
1	A	86	U
1	A	92	U
1	A	93	A
1	A	115	A
1	A	120	A
1	A	124	A
1	A	125	A
1	A	126	A
1	A	129	G
1	A	143	G
1	A	148	G
1	A	150	C
1	A	164	A
1	A	167	C
1	A	182	U
1	A	187	C
1	A	203	G
1	A	212	A
1	A	231	A
1	A	245	C
1	A	246	A
1	A	269	A
1	A	273	A
1	A	283	G
1	A	289	A
1	A	291	A
1	A	292	C
1	A	293	G

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Mol	Chain	Res	Type
1	A	311	U
1	A	314	A
1	A	348	A
1	A	349	U
1	A	368	A
1	A	377	G
1	A	389	A
1	A	397	U
1	A	418	A
1	A	426	U
1	A	427	A
1	A	429	A
1	A	430	U
1	A	431	A
1	A	452	A
1	A	454	C
1	A	456	U
1	A	458	A
1	A	462	A
1	A	467	A
1	A	468	U
1	A	469	U
1	A	470	U
1	A	471	A
1	A	476	A
1	A	488	G
1	A	501	C
1	A	502	A
1	A	503	U
1	A	504	A
1	A	505	G
1	A	515	C
1	A	524	U
1	A	526	G
1	A	529	G
1	A	530	U
1	A	531	A
1	A	532	C
1	A	537	U
1	A	538	A
1	A	539	U
1	A	540	C

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Mol	Chain	Res	Type
1	A	565	U
1	A	570	A
1	A	572	A
1	A	573	C
1	A	603	U
1	A	624	U
1	A	625	C
1	A	629	A
1	A	638	A
1	A	639	A
1	A	643	U
1	A	644	A
1	A	648	A
1	A	655	C
1	A	664	C
1	A	667	A
1	A	680	A
1	A	681	G
1	A	697	A
1	A	709	G
1	A	710	A
1	A	731	A
1	A	742	A
1	A	757	A
1	A	760	C
1	A	767	A
1	A	769	U
1	A	770	A
1	A	776	G
1	A	784	A
1	A	786	U
1	A	797	U
1	A	817	A
1	A	828	A
1	A	832	A
1	A	835	C
1	A	851	C
1	A	860	A
1	A	862	A
1	A	863	U
1	A	864	U
1	A	865	A

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Mol	Chain	Res	Type
1	A	866	A
1	A	867	A
1	A	868	C
1	A	873	C
1	A	874	A
1	A	878	U
1	A	879	U
1	A	880	A
1	A	886	U
1	A	895	G
1	A	911	A
1	A	912	A
1	A	934	U
1	A	948	G
1	A	949	G
1	A	952	U

There are no RNA pucker outliers to report.

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

2 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	$\# Z > 2$
9	5F0	I	181	9	8,8,9	0.57	0	8,9,11	0.99	1 (12%)
17	AYA	Q	2	17	6,7,8	0.78	0	6,8,10	0.50	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
9	5F0	I	181	9	-	0/9/9/10	-
17	AYA	Q	2	17	-	1/5/6/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
9	I	181	5F0	O-C-CB	-2.42	118.34	125.38

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
17	Q	2	AYA	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 45 ligands modelled in this entry, 42 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	$\# Z > 2$
36	FES	P	201	5,16	0,4,4	-	-	-		
37	ATP	X	401	33	32,33,33	0.53	0	48,52,52	0.54	0
36	FES	T	201	13,20	0,4,4	-	-	-		

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	FES	P	201	5,16	-	-	0/1/1/1
37	ATP	X	401	33	-	0/22/38/38	0/3/3/3
36	FES	T	201	13,20	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

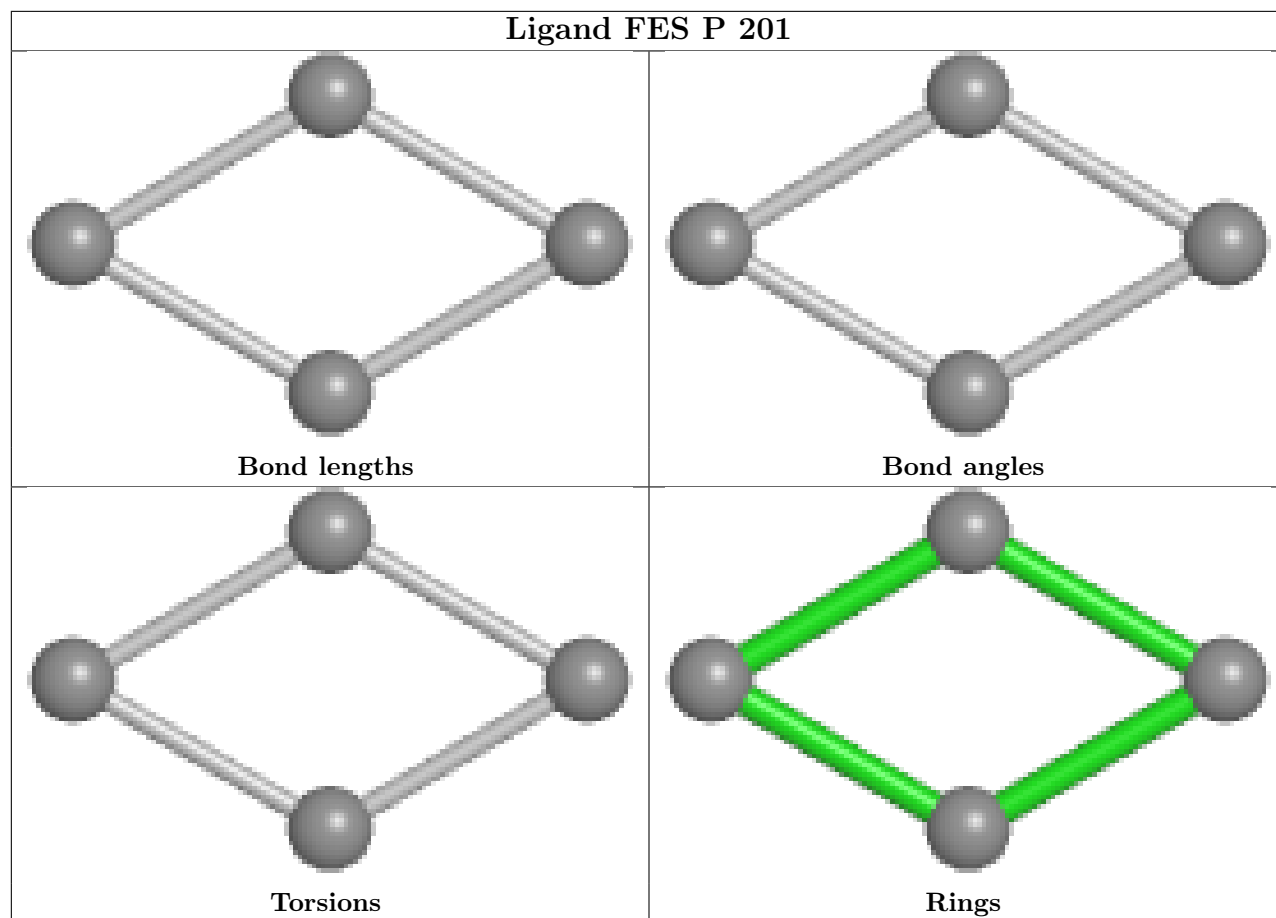
There are no chirality outliers.

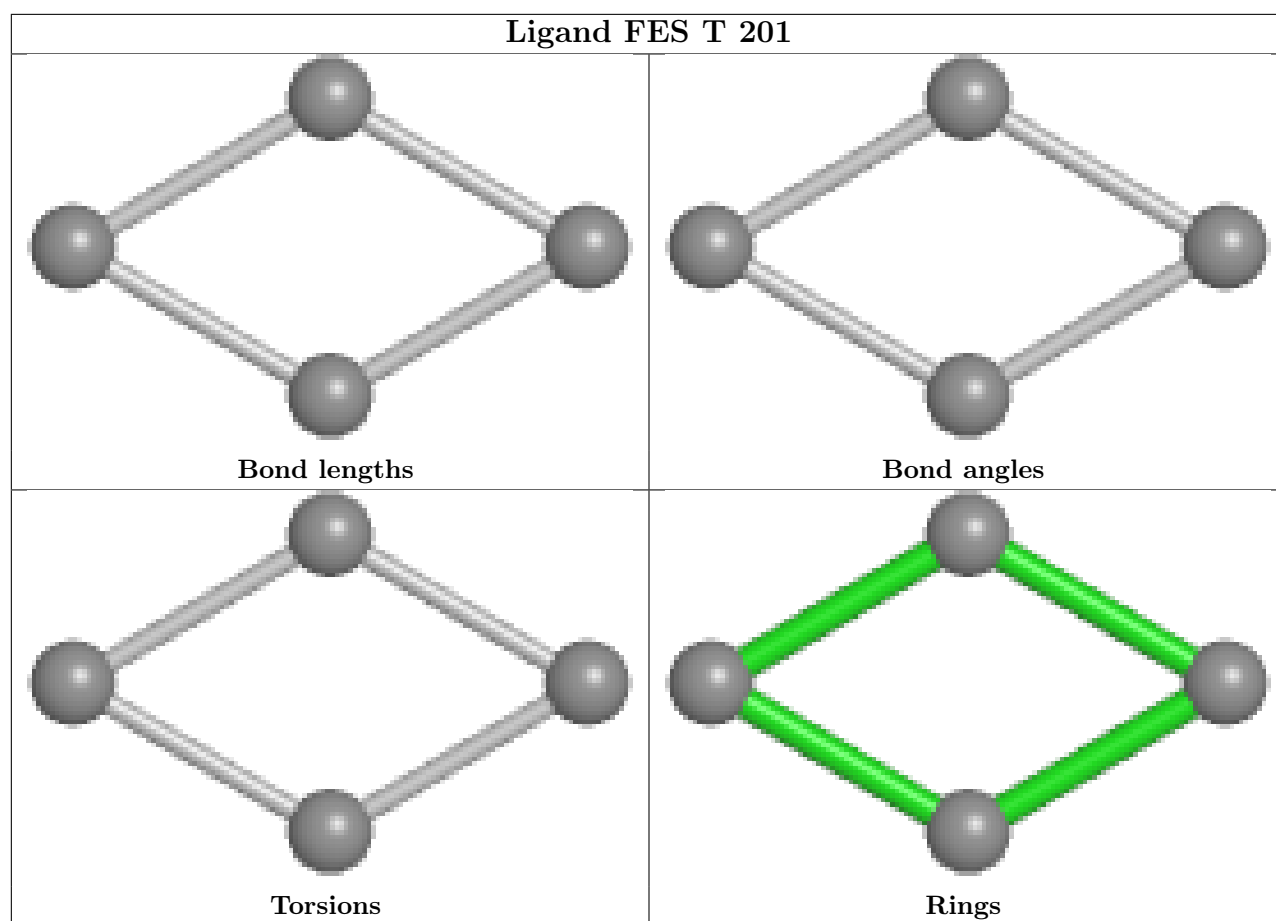
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

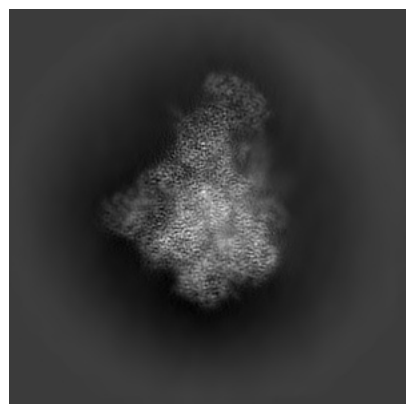
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13551. These allow visual inspection of the internal detail of the map and identification of artifacts.

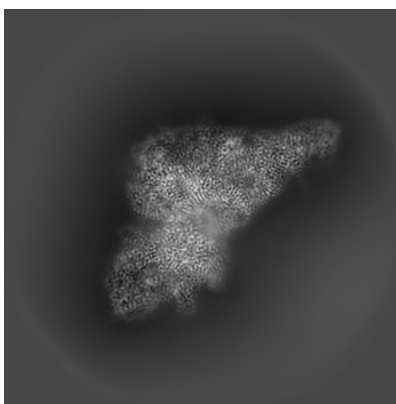
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

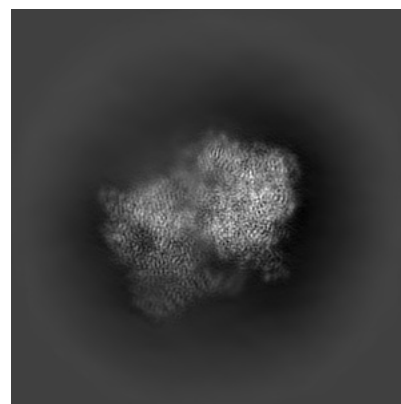
6.1.1 Primary map



X

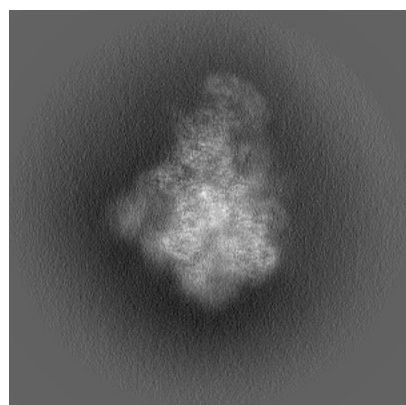


Y

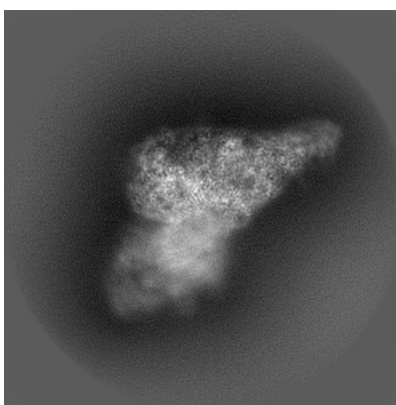


Z

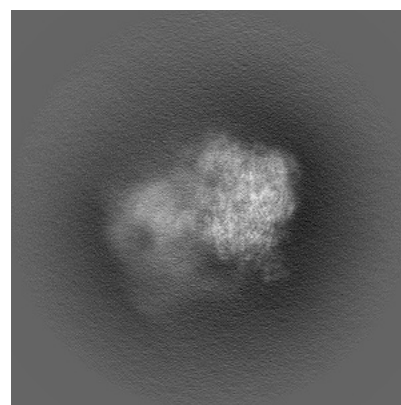
6.1.2 Raw map



X



Y

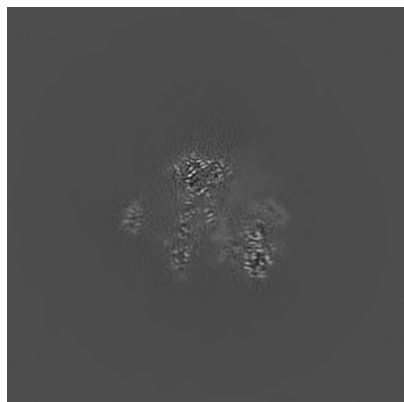


Z

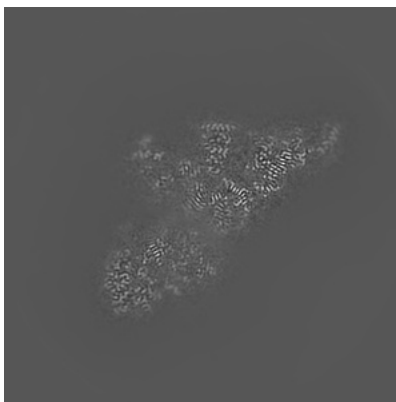
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

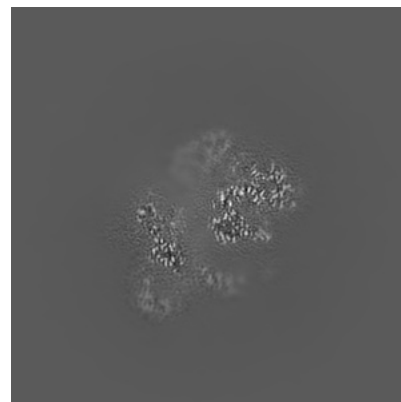
6.2.1 Primary map



X Index: 288

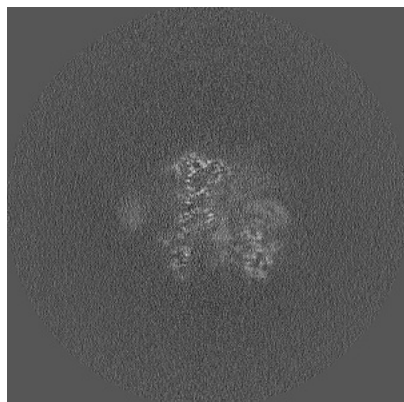


Y Index: 288

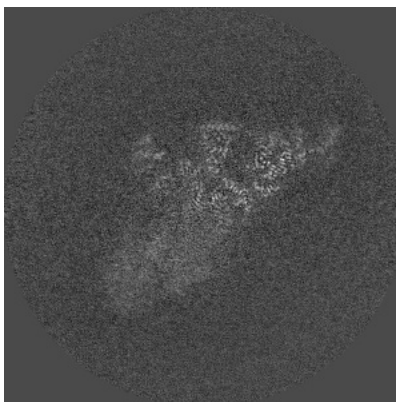


Z Index: 288

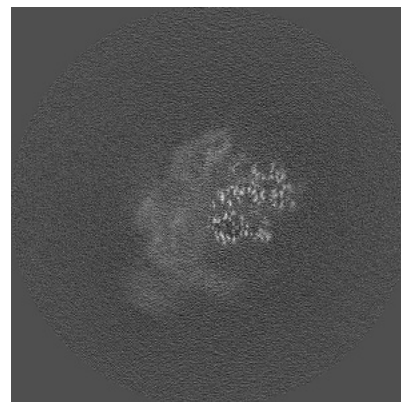
6.2.2 Raw map



X Index: 240



Y Index: 240

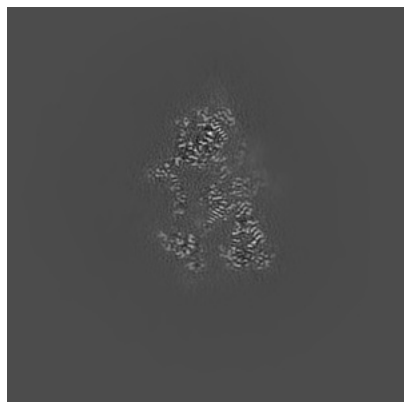


Z Index: 240

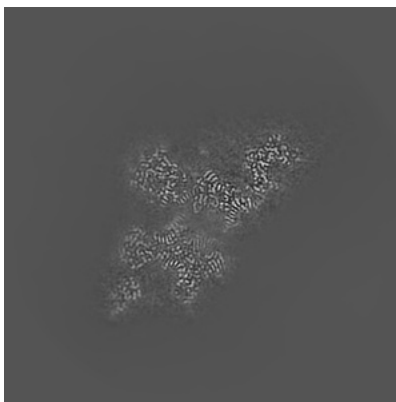
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

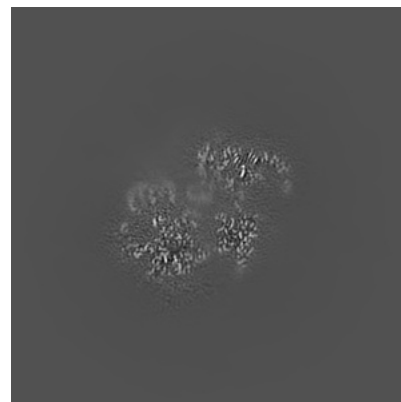
6.3.1 Primary map



X Index: 348

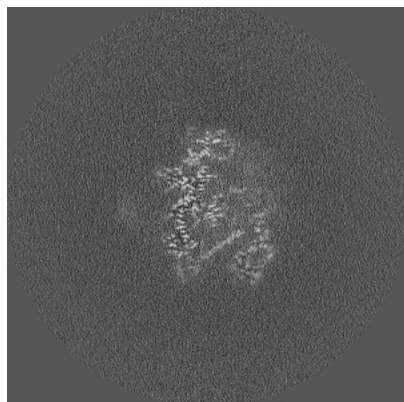


Y Index: 265

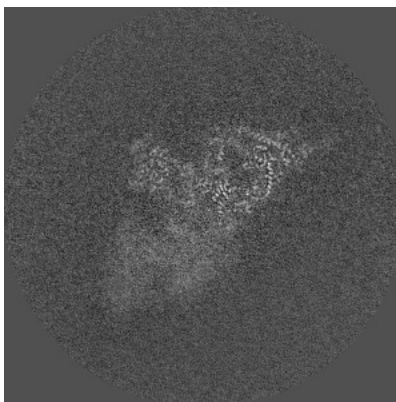


Z Index: 241

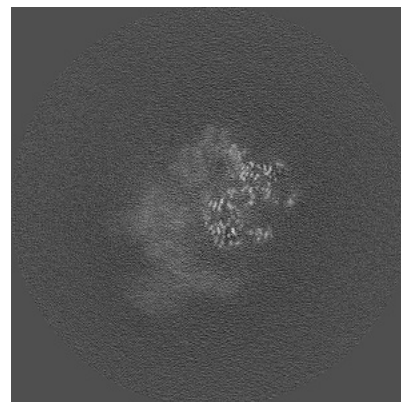
6.3.2 Raw map



X Index: 262



Y Index: 232

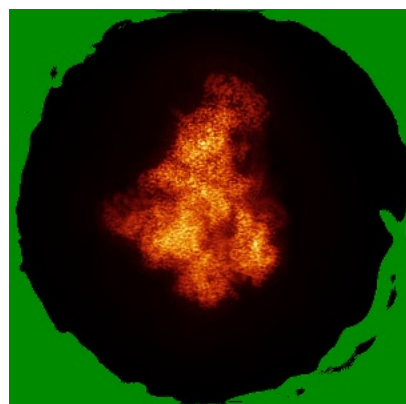


Z Index: 231

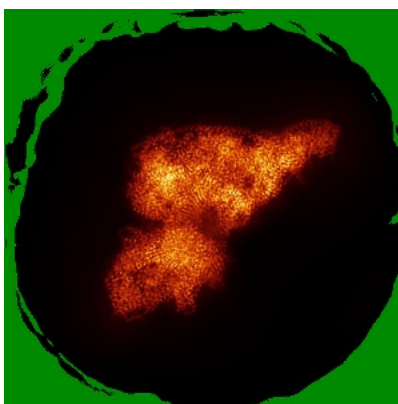
The images above show the largest variance slices of the map in three orthogonal directions.

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

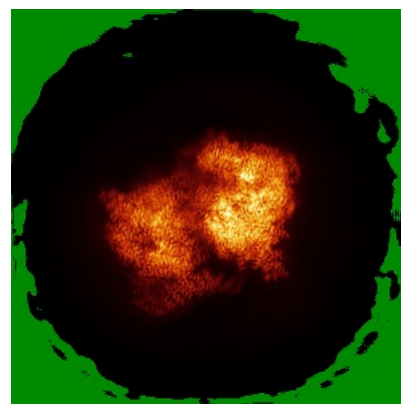
6.4.1 Primary map



X

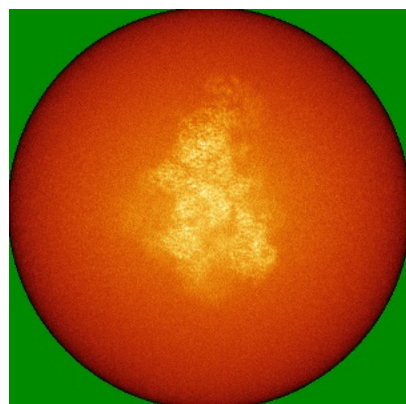


Y

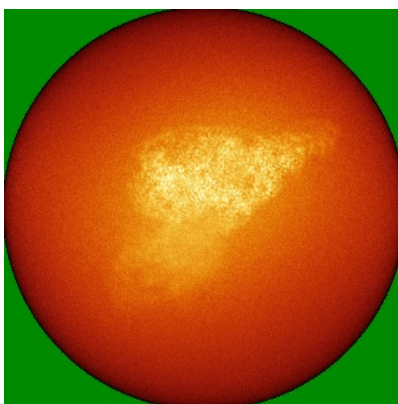


Z

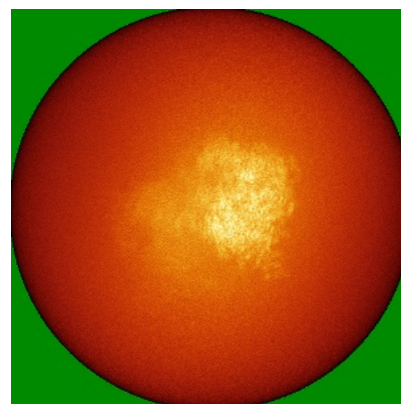
6.4.2 Raw map



X



Y

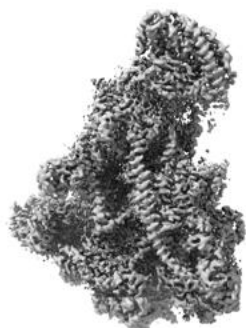


Z

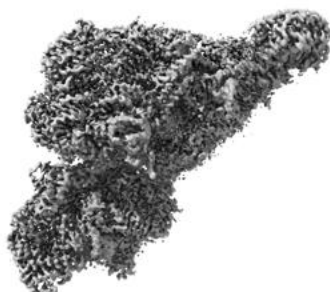
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



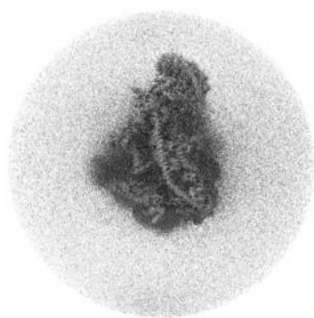
Y



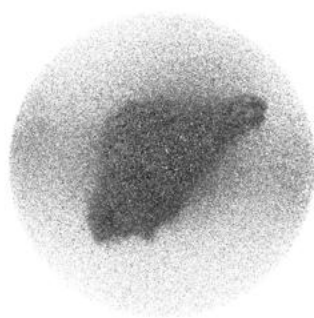
Z

The images above show the 3D surface view of the map at the recommended contour level 0.12. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

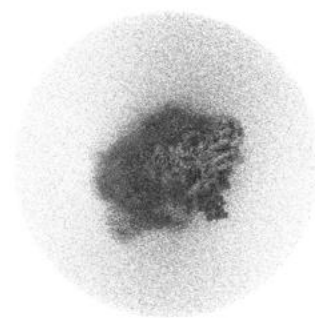
6.5.2 Raw map



X



Y



Z

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

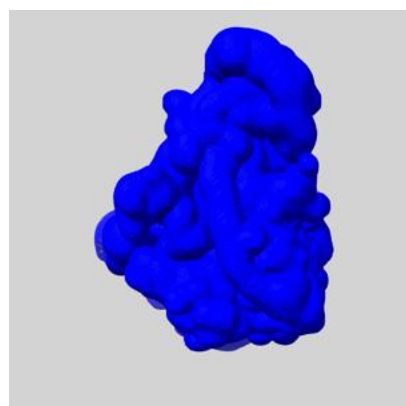
6.6 Mask visualisation [i](#)

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

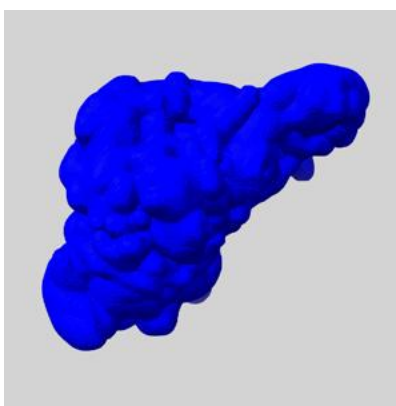
A mask typically either:

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

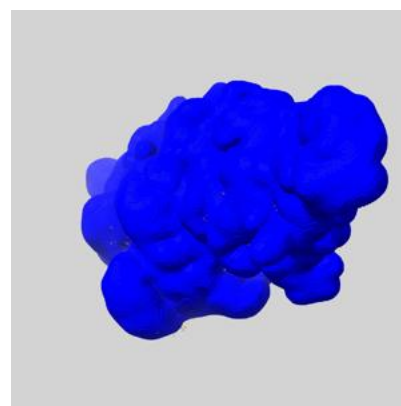
6.6.1 emd_13551_msk_1.map [i](#)



X



Y

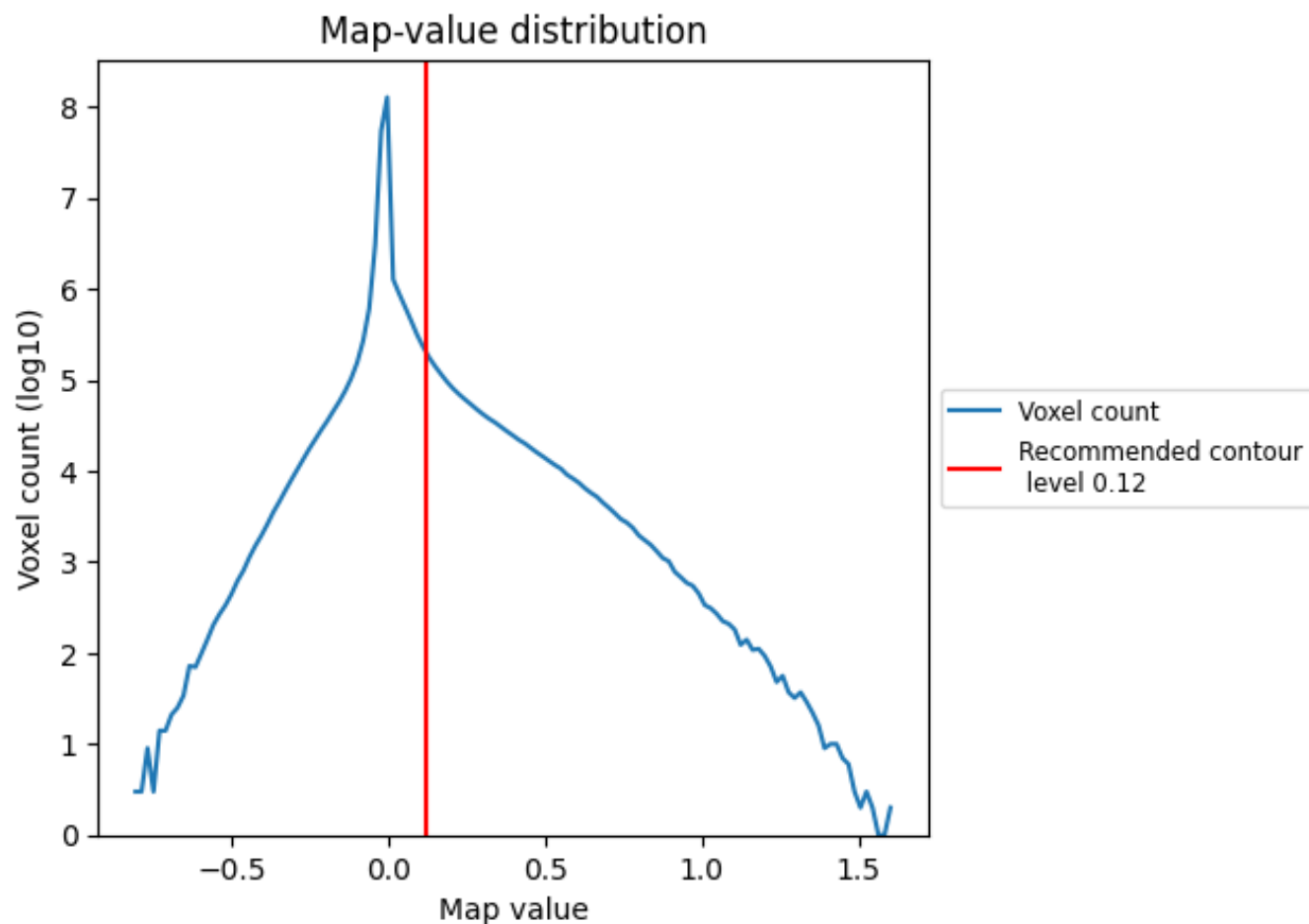


Z

7 Map analysis [i](#)

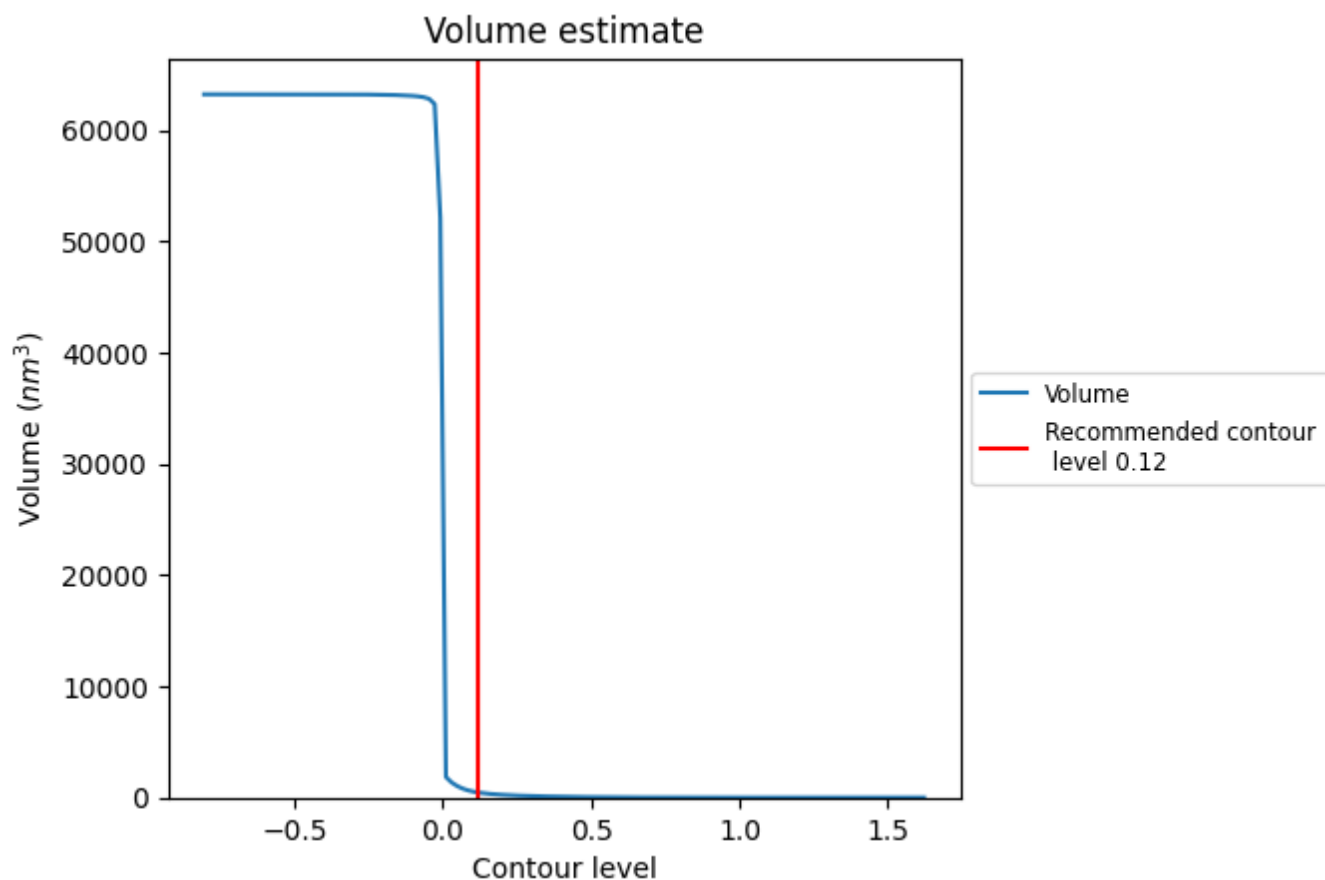
This section contains the results of statistical analysis of the map.

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.

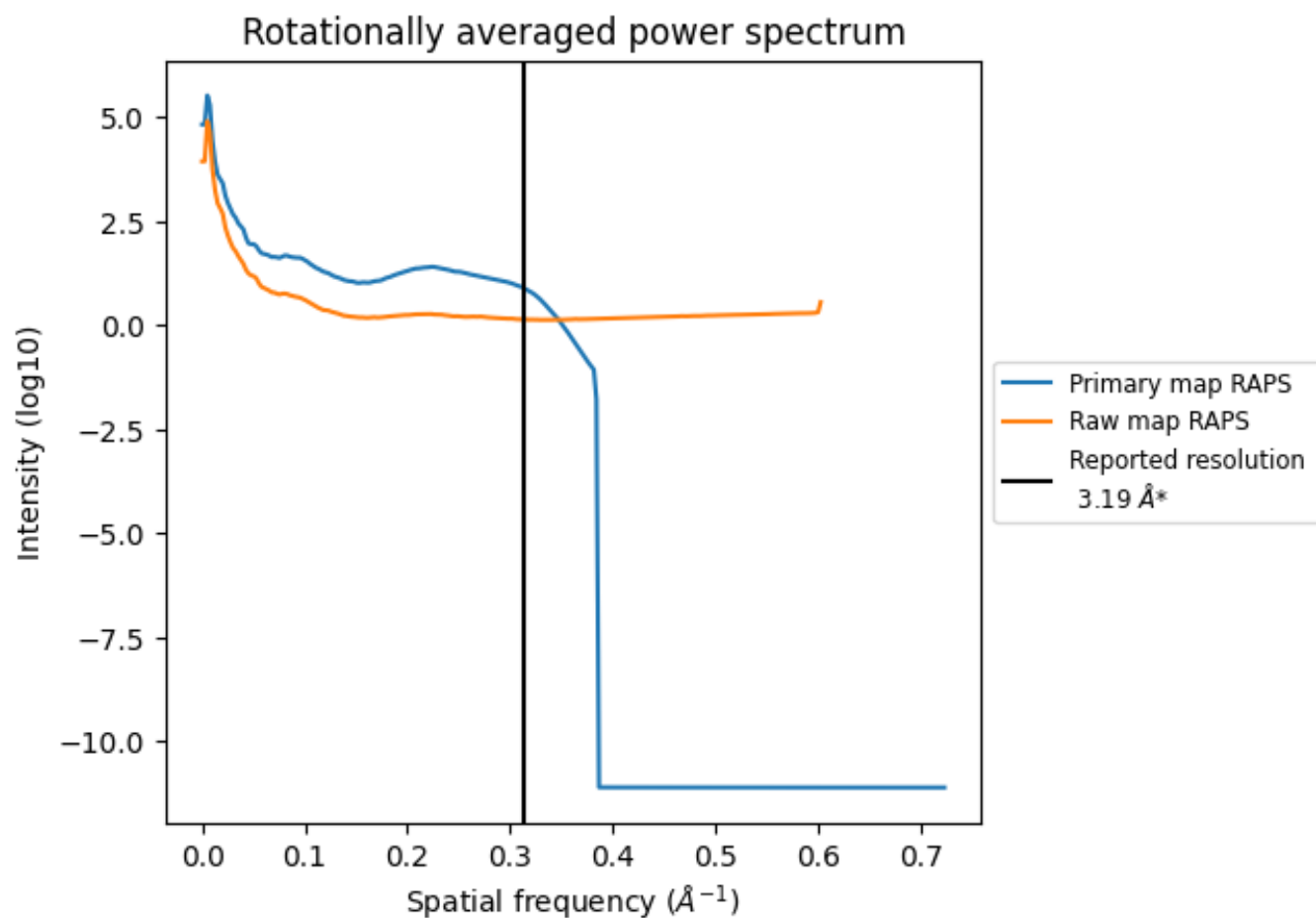
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 453 nm³; this corresponds to an approximate mass of 410 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 ⓘ

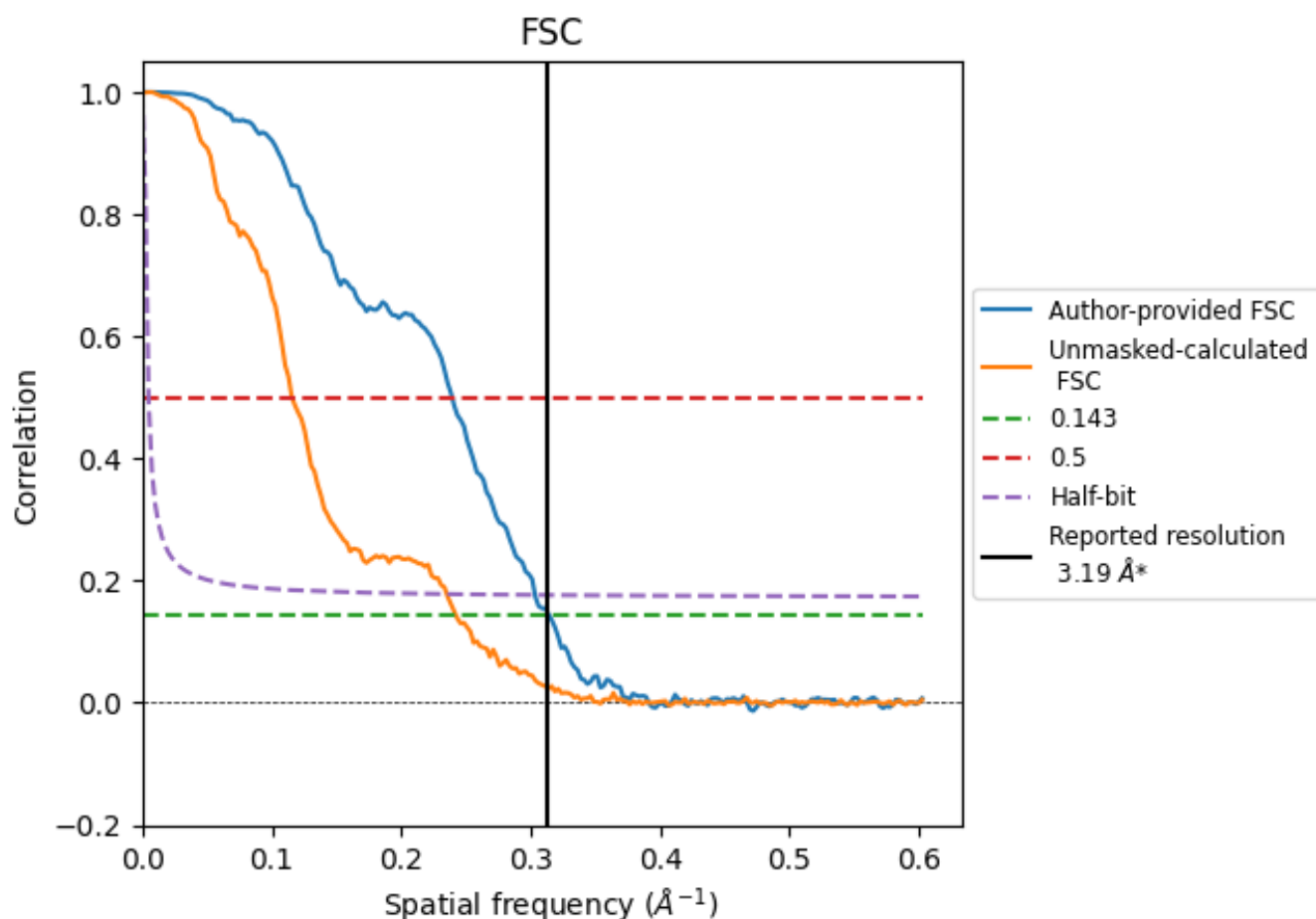


*Reported resolution corresponds to spatial frequency of 0.313 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.313 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

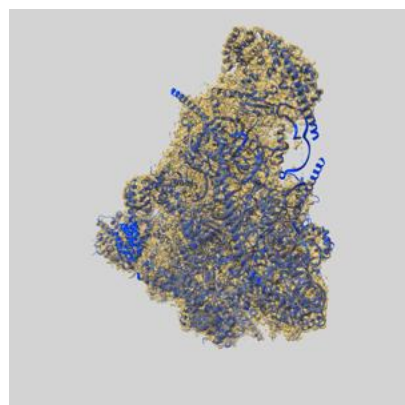
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.19	-	-
Author-provided FSC curve	3.18	4.18	3.30
Unmasked-calculated*	4.13	8.66	4.25

*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 4.13 differs from the reported value 3.19 by more than 10 %

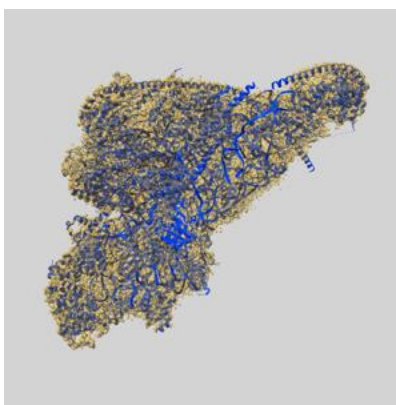
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13551 and PDB model 7PNT. Per-residue inclusion information can be found in section 3 on page 13.

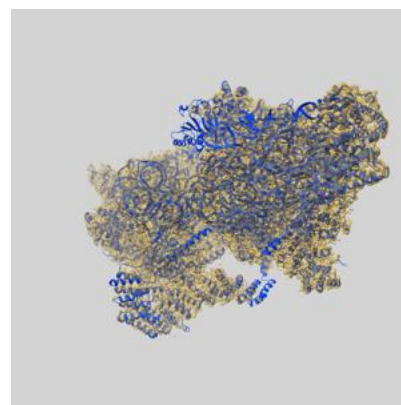
9.1 Map-model overlay [i](#)



X



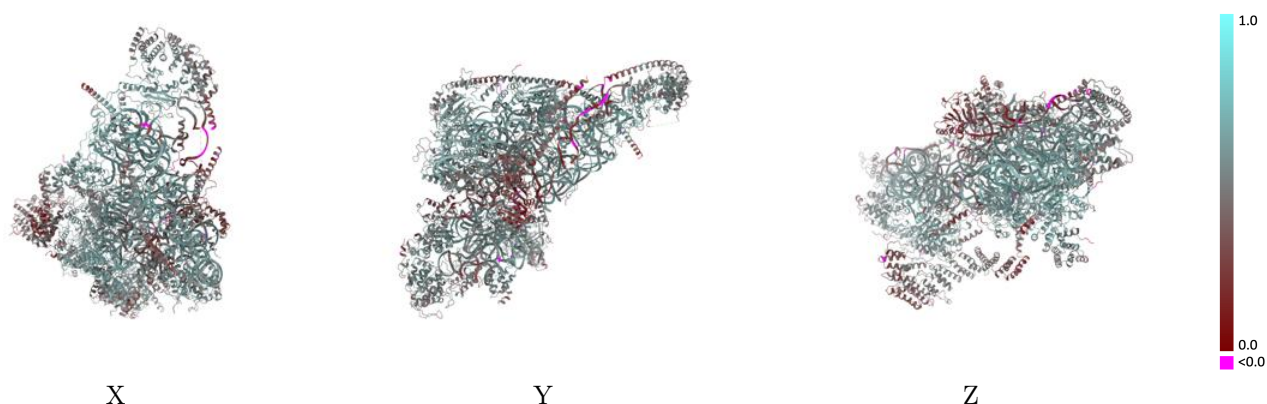
Y



Z

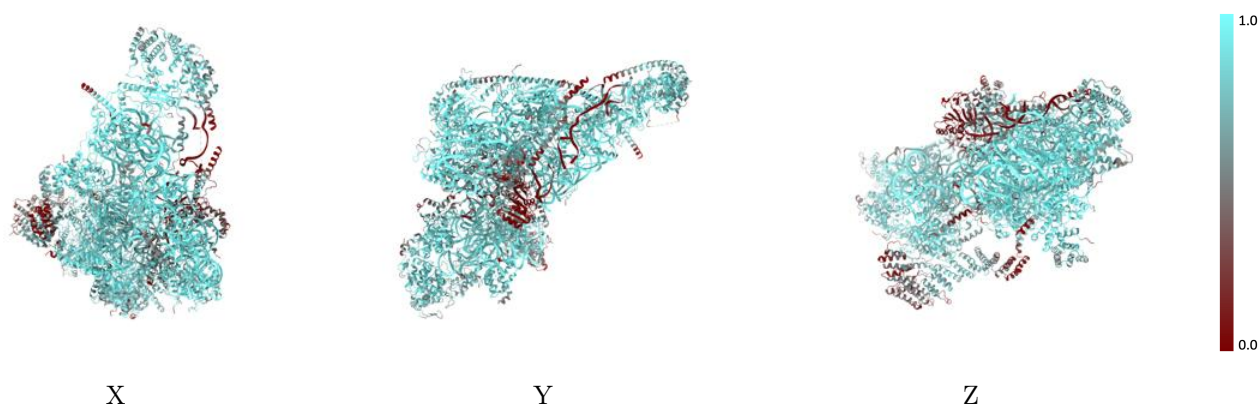
The images above show the 3D surface view of the map at the recommended contour level 0.12 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



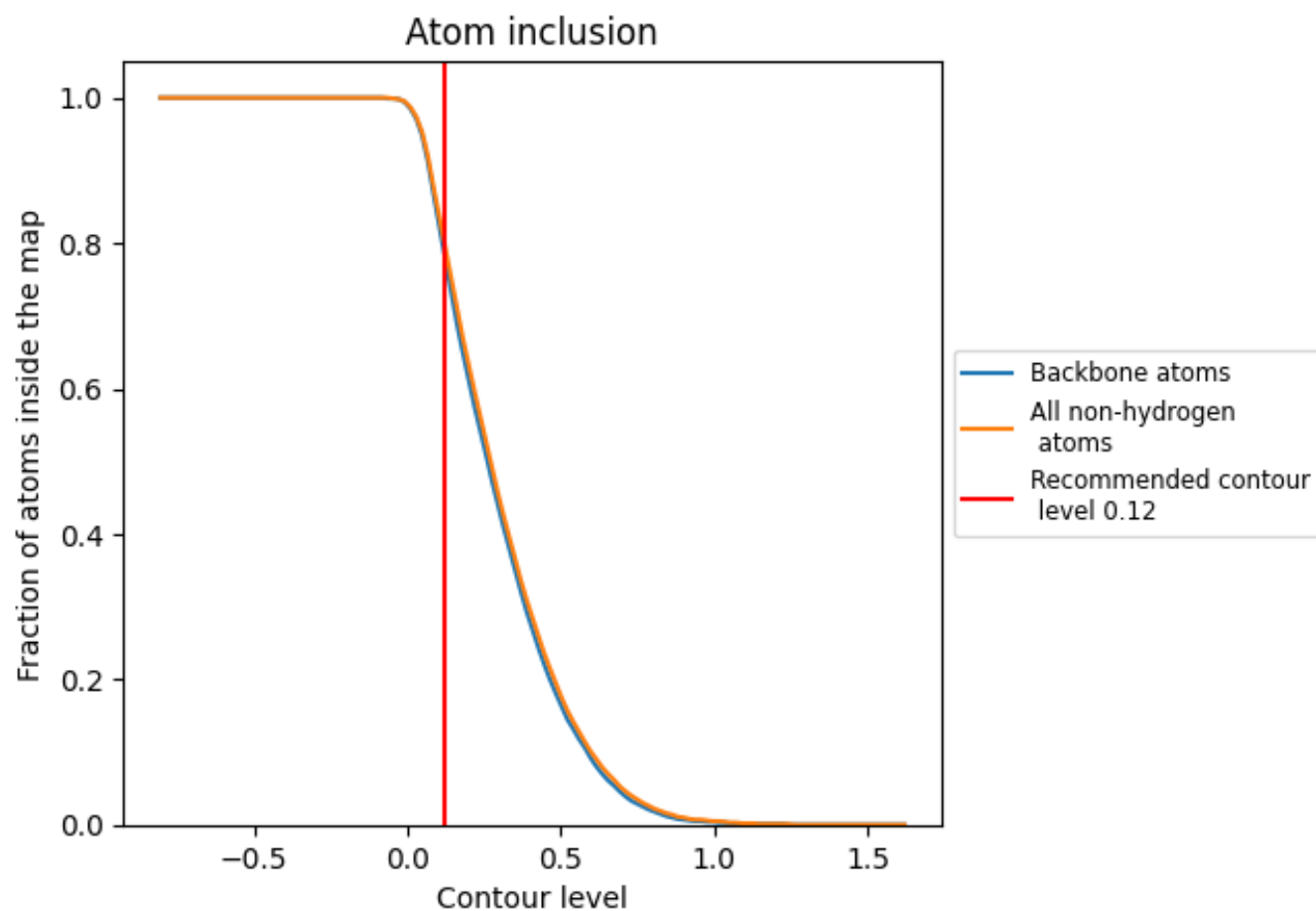
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.12).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 80% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8020	 0.5160
0	 0.8830	 0.5530
1	 0.8110	 0.5240
3	 0.2940	 0.3520
4	 0.5530	 0.4090
A	 0.8700	 0.5360
B	 0.9280	 0.6050
C	 0.9350	 0.5990
D	 0.8100	 0.5270
E	 0.8750	 0.5470
F	 0.7720	 0.4890
G	 0.7910	 0.5150
H	 0.8730	 0.5570
I	 0.9060	 0.5570
J	 0.8870	 0.5300
K	 0.9370	 0.5950
L	 0.7720	 0.5040
M	 0.8940	 0.6000
N	 0.8970	 0.5800
O	 0.9250	 0.5970
P	 0.9140	 0.5750
Q	 0.9090	 0.5620
R	 0.8540	 0.5550
S	 0.8020	 0.5040
T	 0.9020	 0.5750
U	 0.7710	 0.5010
V	 0.7410	 0.4880
W	 0.8710	 0.5640
X	 0.8140	 0.5200
Y	 0.6550	 0.4620
Z	 0.7990	 0.5180
a	 0.5390	 0.4140
c	 0.2460	 0.2710

