



# wwPDB EM Validation Summary Report ⓘ

Dec 29, 2024 – 10:22 AM EST

PDB ID : 7OF0  
EMDB ID : EMD-12865  
Title : Structure of a human mitochondrial ribosome large subunit assembly intermediate in complex with MTERF4-NSUN4 (dataset1).  
Authors : Hillen, H.S.; Lavdovskaia, E.; Nadler, F.; Hanitsch, E.; Linden, A.; Bohnsack, K.E.; Urlaub, H.; Richter-Dennerlein, R.  
Deposited on : 2021-05-04  
Resolution : 2.20 Å (reported)  
Based on initial model : 5OOL

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev113  
MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.40

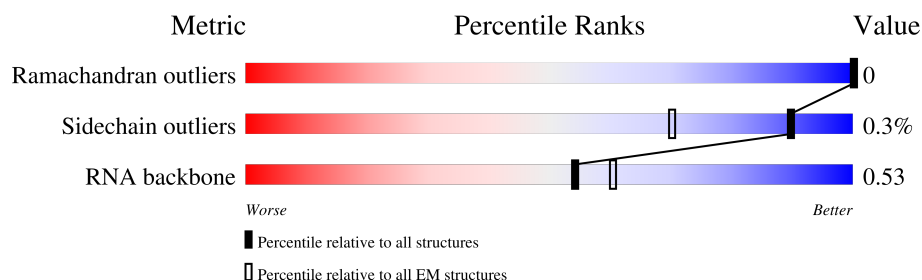
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.20 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.















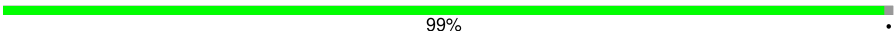

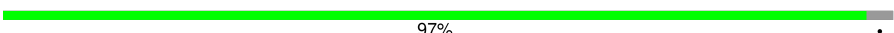







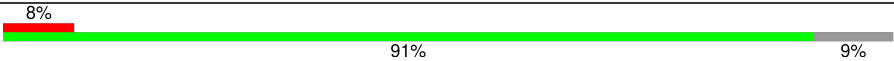


Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	57% 43%
2	1	65	80% 20%
3	2	92	49% 51%
4	3	188	51% 49%
5	4	103	36% 64%
6	5	423	92% 7%
7	6	380	85% 15%
8	7	338	83% 15%

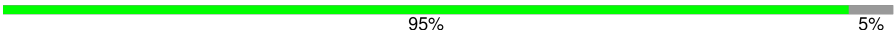










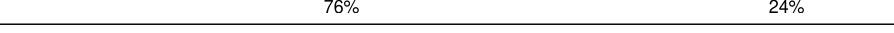
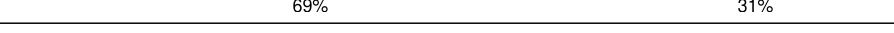
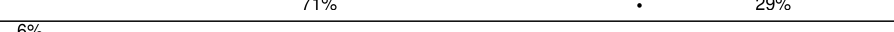

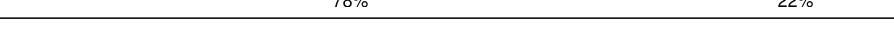



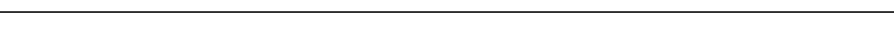


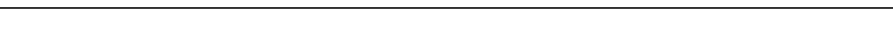
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Mol	Chain	Length	Quality of chain
9	8	206	
10	9	137	
11	A	1559	
12	B	69	
13	C	384	
14	D	305	
15	E	348	
16	F	311	
17	G	381	
18	H	267	
19	I	261	
20	J	192	
21	K	178	
22	L	145	
23	M	296	
24	N	251	
25	O	175	
26	P	180	
27	Q	292	
28	R	149	
29	S	205	
30	T	206	
31	U	153	
32	V	216	
33	W	148	

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Mol	Chain	Length	Quality of chain
34	X	256	 95% 5%
35	Y	250	 70% 30%
36	Z	161	 71% 29%
37	a	142	 50% 50%
38	b	215	 69% 31%
39	c	332	 83% 17%
40	d	306	 23% 64% 35%
41	e	279	 9% 71% 29%
42	f	212	 9% 51% 49%
43	g	166	 77% 22%
44	h	158	 66% 34%
45	i	128	 76% 24%
46	j	123	 69% 31%
47	k	112	 6% 71% 29%
48	m	128	 6% 23% 77%
49	o	102	 78% 22%
50	p	206	 62% 38%
51	q	222	 58% 42%
52	r	196	 74% 26%
53	s	439	 84% 16%
54	u	234	 47% 53%
55	v	70	 96% ..
56	w	156	 50% 49%

## 2 Entry composition

There are 58 unique types of molecules in this entry. The entry contains 99476 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	52	Total	C	N	O	S	0	0
			433	278	83	70	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	45	Total	C	N	O	S	0	0
			367	227	81	58	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	37	Total	C	N	O	S	0	0
			333	212	71	47	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
6	5	392	Total	C	N	O	S	0	0
			3199	2067	558	563	11		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
7	6	324	Total	C	N	O	S	0	0
			2723	1743	488	484	8		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
8	7	287	Total	C	N	O	S	0	0
			2334	1495	397	425	17		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
9	8	77	Total	C	N	O	S	0	0
			651	413	113	123	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
10	9	117	Total	C	N	O	S	0	0
			947	614	163	168	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
11	A	1376	Total	C	N	O	P	0	0
			29219	13111	5277	9455	1376		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3107	U	UNK	conflict	GB 1025814679

- Molecule 12 is a RNA chain called mitochondrial tRNA<sup>Val</sup>.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	B	56	Total	C	N	O	P	0	0
			1191	534	214	387	56		

- Molecule 13 is a protein called 5-methylcytosine rRNA methyltransferase NSUN4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	C	327	Total	C	N	O	S	1	0
			2595	1651	454	474	16		

- Molecule 14 is a protein called 39S ribosomal protein L2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	D	240	Total	C	N	O	S	0	0
			1872	1165	378	320	9		

- Molecule 15 is a protein called 39S ribosomal protein L3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	E	308	Total	C	N	O	S	0	0
			2427	1559	423	434	11		

- Molecule 16 is a protein called 39S ribosomal protein L4, mitochondrial.

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

- Molecule 17 is a protein called Transcription termination factor 4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	G	238	Total	C	N	O	S	0	0
			1943	1243	336	352	12		

- Molecule 18 is a protein called 39S ribosomal protein L9, mitochondrial.

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

- Molecule 19 is a protein called 39S ribosomal protein L10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	I	151	Total	C	N	O	S	0	0
			1228	795	222	201	10		

- Molecule 20 is a protein called 39S ribosomal protein L11, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	J	140	Total	C	N	O	S	0	0
			1061	680	192	187	2		

- Molecule 21 is a protein called 39S ribosomal protein L13, mitochondrial.

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

- Molecule 22 is a protein called 39S ribosomal protein L14, mitochondrial.

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

- Molecule 23 is a protein called 39S ribosomal protein L15, mitochondrial.

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

- Molecule 24 is a protein called 39S ribosomal protein L16, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	N	183	Total	C	N	O	S	0	0
			1494	960	274	251	9		

- Molecule 25 is a protein called 39S ribosomal protein L17, mitochondrial.

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

- Molecule 26 is a protein called 39S ribosomal protein L18, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	P	141	Total	C	N	O	S	0	0
			1148	719	221	203	5		

- Molecule 27 is a protein called 39S ribosomal protein L19, mitochondrial.



Mol	Chain	Residues	Atoms					AltConf	Trace
27	Q	217	Total	C	N	O	S	0	0
			1805	1159	317	320	9		

- Molecule 28 is a protein called 39S ribosomal protein L20, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	R	116	Total	C	N	O	S	0	0
			971	616	200	151	4		

- Molecule 29 is a protein called 39S ribosomal protein L21, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	S	156	Total	C	N	O	S	0	0
			1251	806	222	219	4		

- Molecule 30 is a protein called 39S ribosomal protein L22, mitochondrial.

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

- Molecule 31 is a protein called 39S ribosomal protein L23, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	U	139	Total	C	N	O	S	0	0
			1154	734	220	197	3		

- Molecule 32 is a protein called 39S ribosomal protein L24, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	V	47	Total	C	N	O	S	0	0
			395	255	62	75	3		

- Molecule 33 is a protein called 39S ribosomal protein L27, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	W	100	Total	C	N	O	S	0	0
			801	518	150	130	3		

- Molecule 34 is a protein called 39S ribosomal protein L28, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	X	243	Total	C	N	O	S	0	0
			2035	1317	351	362	5		

- Molecule 35 is a protein called 39S ribosomal protein L47, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	Y	176	Total	C	N	O	S	0	0
			1517	970	291	252	4		

- Molecule 36 is a protein called 39S ribosomal protein L30, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	Z	115	Total	C	N	O	S	0	0
			937	598	175	161	3		

- Molecule 37 is a protein called 39S ribosomal protein L42, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	a	71	Total	C	N	O	S	0	0
			597	378	112	103	4		

- Molecule 38 is a protein called 39S ribosomal protein L43, mitochondrial.

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

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

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

- Molecule 40 is a protein called 39S ribosomal protein L45, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	d	199	Total	C	N	O	S	0	0
			1653	1075	276	293	9		

- Molecule 41 is a protein called 39S ribosomal protein L46, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	e	197	Total	C	N	O	S	0	0
			1599	1027	277	290	5		

- Molecule 42 is a protein called 39S ribosomal protein L48, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	f	108	Total	C	N	O	S	0	0
			857	549	140	165	3		

- Molecule 43 is a protein called 39S ribosomal protein L49, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	g	129	Total	C	N	O	S	0	0
			1067	690	185	190	2		

- Molecule 44 is a protein called 39S ribosomal protein L50, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	h	105	Total	C	N	O	S	0	0
			862	548	151	160	3		

- Molecule 45 is a protein called 39S ribosomal protein L51, mitochondrial.

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

- Molecule 46 is a protein called 39S ribosomal protein L52, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	j	85	Total	C	N	O	S	0	0
			684	423	133	126	2		

- Molecule 47 is a protein called 39S ribosomal protein L53, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	k	80	Total	C	N	O	S	0	0
			627	392	116	114	5		

- Molecule 48 is a protein called 39S ribosomal protein L55, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	m	29	Total	C	N	O	S	0	0
			245	157	48	38	2		

- Molecule 49 is a protein called Ribosomal protein 63, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	o	80	Total	C	N	O	S	0	0
			676	426	134	113	3		

- Molecule 50 is a protein called Peptidyl-tRNA hydrolase ICT1, mitochondrial.

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

- Molecule 51 is a protein called Growth arrest and DNA damage-inducible proteins-interacting protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	q	128	Total	C	N	O	S	0	0
			1076	671	208	192	5		

- Molecule 52 is a protein called 39S ribosomal protein S18a, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	r	146	Total	C	N	O	S	0	0
			1203	764	232	199	8		

- Molecule 53 is a protein called 39S ribosomal protein S30, mitochondrial.

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

- Molecule 54 is a protein called Mitochondrial assembly of ribosomal large subunit protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	u	111	Total	C	N	O	S	0	0
			927	595	155	167	10		

- Molecule 55 is a protein called MIEF1 upstream open reading frame protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
55	v	69	Total	C	N	O	0	0
			588	372	116	100		

- Molecule 56 is a protein called Acyl carrier protein, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	w	79	Total	C	N	O	S	0	0
			638	410	95	128	5		

- Molecule 57 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
57	0	1	Total	Zn	0
			1	1	
57	4	1	Total	Zn	0
			1	1	
57	r	1	Total	Zn	0
			1	1	

- Molecule 58 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

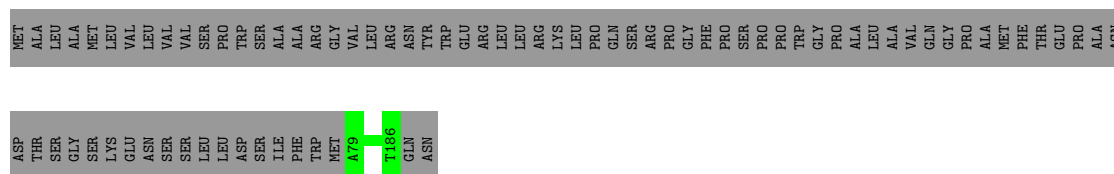
Mol	Chain	Residues	Atoms		AltConf
58	A	59	Total	Mg	0
			59	59	
58	E	1	Total	Mg	0
			1	1	
58	g	1	Total	Mg	0
			1	1	

### 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.

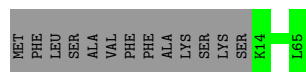
- Molecule 1: 39S ribosomal protein L32, mitochondrial

Chain 0:  57% 43%



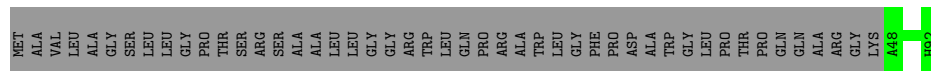
- Molecule 2: 39S ribosomal protein L33, mitochondrial

Chain 1:  80% 20%



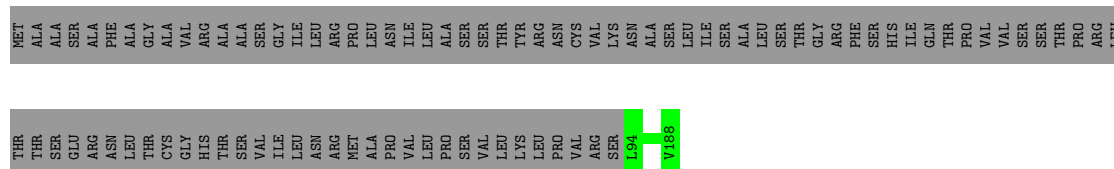
- Molecule 3: 39S ribosomal protein L34, mitochondrial

Chain 2:  49% 51%



- Molecule 4: 39S ribosomal protein L35, mitochondrial

Chain 3:  51% 49%



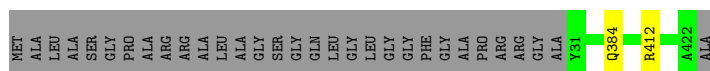
- Molecule 5: 39S ribosomal protein L36, mitochondrial

Chain 4:  36% 64%




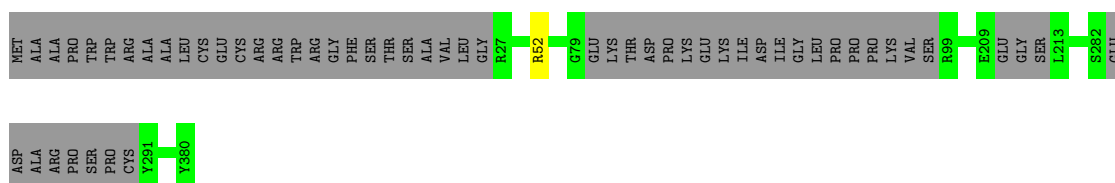
- Molecule 6: 39S ribosomal protein L37, mitochondrial

Chain 5:  92% 7%




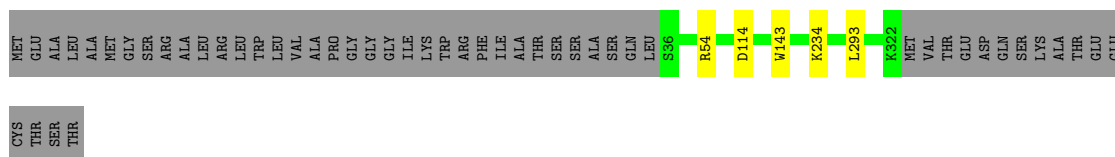
- Molecule 7: 39S ribosomal protein L38, mitochondrial

Chain 6:  85% 15%



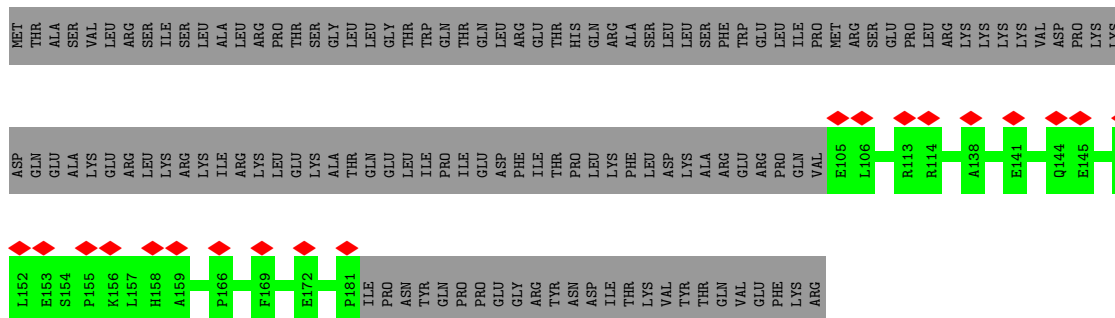
- Molecule 8: 39S ribosomal protein L39, mitochondrial

Chain 7:  83% 15%




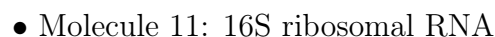
- Molecule 9: 39S ribosomal protein L40, mitochondrial

Chain 8:  10% 37% 63%



- Molecule 10: 39S ribosomal protein L41, mitochondrial

Chain 9:  85% 15%



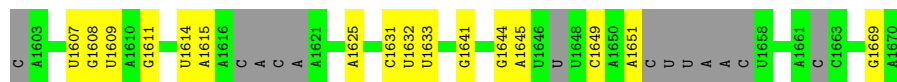
Response	Percentage
Yes	67%
No	21%
Don't know	12%






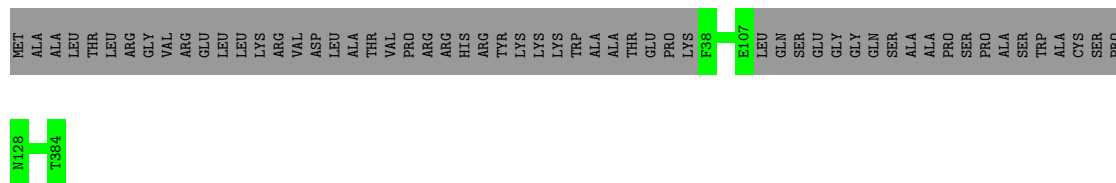
- Molecule 12: mitochondrial tRNA<sup>Val</sup>

Chain B:  58% 23% 19%




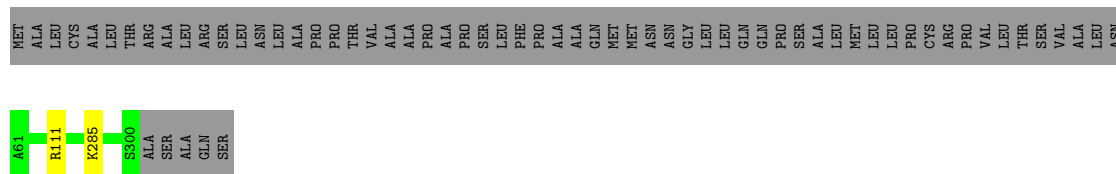
- Molecule 13: 5-methylcytosine rRNA methyltransferase NSUN4

Chain C:  85% 15%



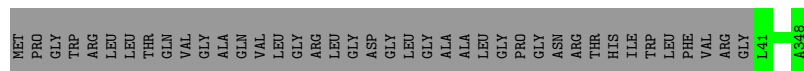
- Molecule 14: 39S ribosomal protein L2, mitochondrial

Chain D:  78% 21%




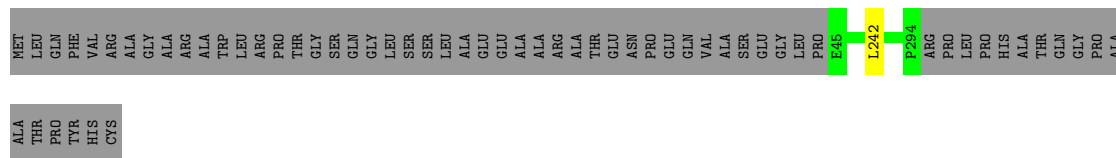
- Molecule 15: 39S ribosomal protein L3, mitochondrial

Chain E:  89% 11%



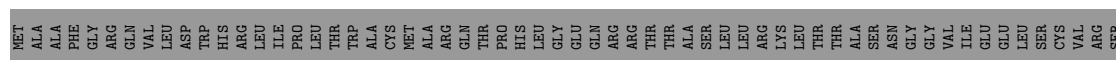
- Molecule 16: 39S ribosomal protein L4, mitochondrial

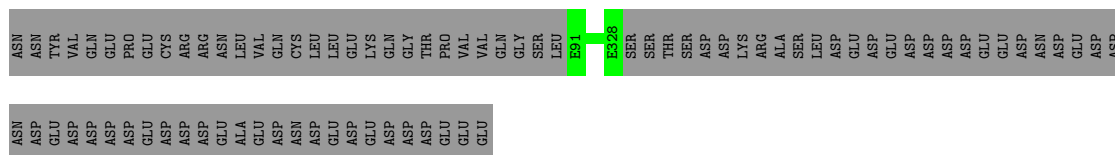
Chain F:  80% 20%



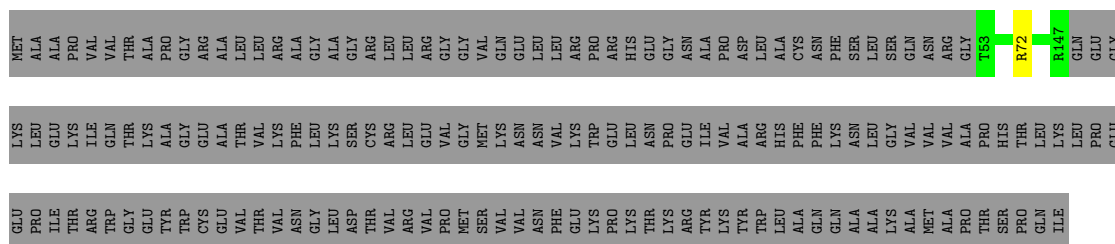
- Molecule 17: Transcription termination factor 4, mitochondrial

Chain G:  62% 38%

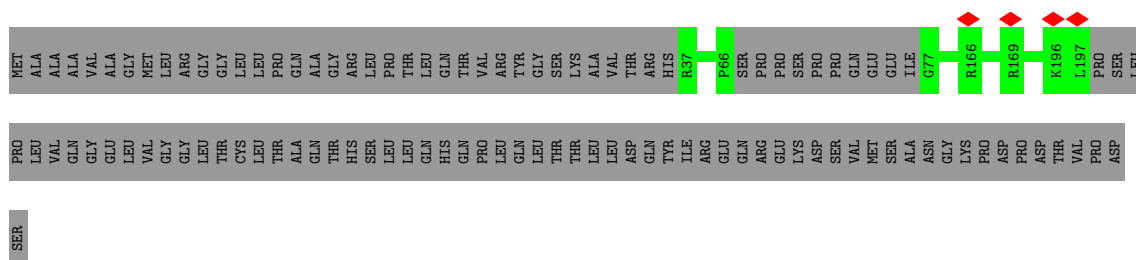




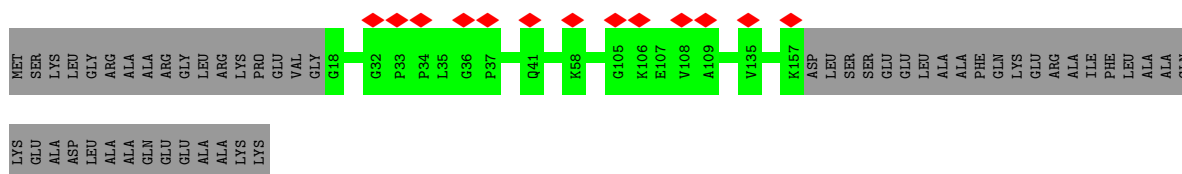
- Molecule 18: 39S ribosomal protein L9, mitochondrial



- Molecule 19: 39S ribosomal protein L10, mitochondrial



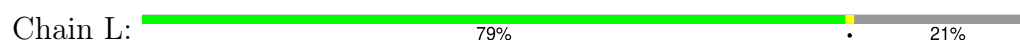
- Molecule 20: 39S ribosomal protein L11, mitochondrial

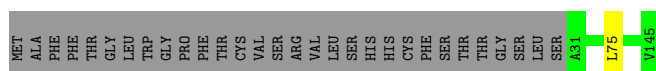


- Molecule 21: 39S ribosomal protein L13, mitochondrial



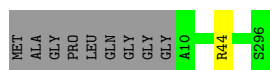
- Molecule 22: 39S ribosomal protein L14, mitochondrial





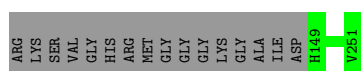
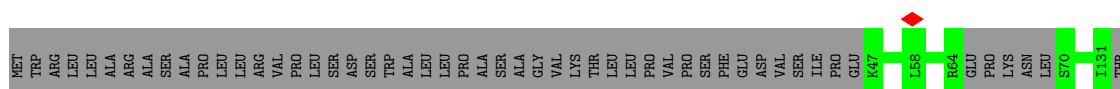
- Molecule 23: 39S ribosomal protein L15, mitochondrial

Chain M: 97%



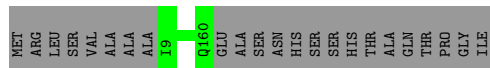
- Molecule 24: 39S ribosomal protein L16, mitochondrial

Chain N: 73% 27%



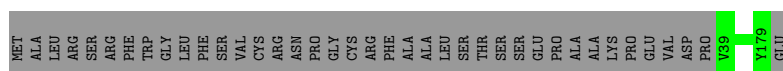
- Molecule 25: 39S ribosomal protein L17, mitochondrial

Chain O: 87% 13%



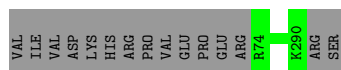
- Molecule 26: 39S ribosomal protein L18, mitochondrial

Chain P: 78% 22%



- Molecule 27: 39S ribosomal protein L19, mitochondrial

Chain Q: 74% 26%



- Molecule 28: 39S ribosomal protein L20, mitochondrial

Chain R: 78% 22%



- Molecule 35: 39S ribosomal protein L47, mitochondrial

Chain Y:  70% 30%

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

ARG LYS **G63** **F238** PRO HIS LEU ALA GLU ALA GLN LYS SER SER SER LEU VAL

- Molecule 36: 39S ribosomal protein L30, mitochondrial

Chain Z:  71% 29%

MET ALA GLY ILE LEU LEU ARG VAL VAL VAL GLN TRP PRO PRO GLY ARG GLN THR VAL THR LYS GLY VAL GLN SER LEU ILE CYS THR ASP TRP ILE ARG HIS **K35** **K44** **W149** HIS LEU LYS PRO VAL GLU GLN LYS ALA HIS GLU SER

- Molecule 37: 39S ribosomal protein L42, mitochondrial

Chain a:  50% 50%

MET VAL ALA VAL ALA ALA TRP VAL MET MET LYS ARG THR ILE LEU LYS HIS LEU PHE PRO VAL GLN ASN GLY ALA LEU TYR CYS VAL CYS HIS LYS SER THR TYR SER PRO LEU PRO ASP TYR ASN CYS **N46** **R77** PRO ASP PRO VAL HIS ASN ASN GLU THR HIS PRO ASP

GLN VAL LEU LYS THR ARG LEU GLU VAL LYS VAL HIS LEU **E104** **R142**


- Molecule 38: 39S ribosomal protein L43, mitochondrial

Chain b:  69% 31%

MET **T2** **Q149** ASP PRO PRO ALA ALA ALA GLN ASP THR GLY LEU ARG LEU SER ALA VAL ALA PRO PRO VAL ALA PRO PRO ILE ILE LEU LEU PRO PRO GLY TRP ARG GLY **V31** **Q107** LEU LEU ILE GLY GLY LYS GLU VAL SER SER SER LEU ASN **L119** **Y316** SER LYS PRO LYS THR THR LEU ARG ALA GLU LYS PRO HIS PRO ILE

VAL PRO ALA LEU THR VAL CYS SER ALA

- Molecule 39: 39S ribosomal protein L44, mitochondrial

Chain c:  83% 17%

MET ALA SER GLY LEU VAL ARG LEU LEU GLN GLN GLY HIS ARG CYS LEU LEU ALA PRO VAL ALA PRO PRO LYS LEU VAL VAL PRO VAL ARG GLY **V31** **Q107** LEU LEU ILE GLY GLY LYS GLU VAL SER SER SER LEU ASN **L119** **Y316** SER LYS PRO LYS THR THR LEU ARG ALA GLU LYS PRO HIS PRO ILE

THR ALA SER

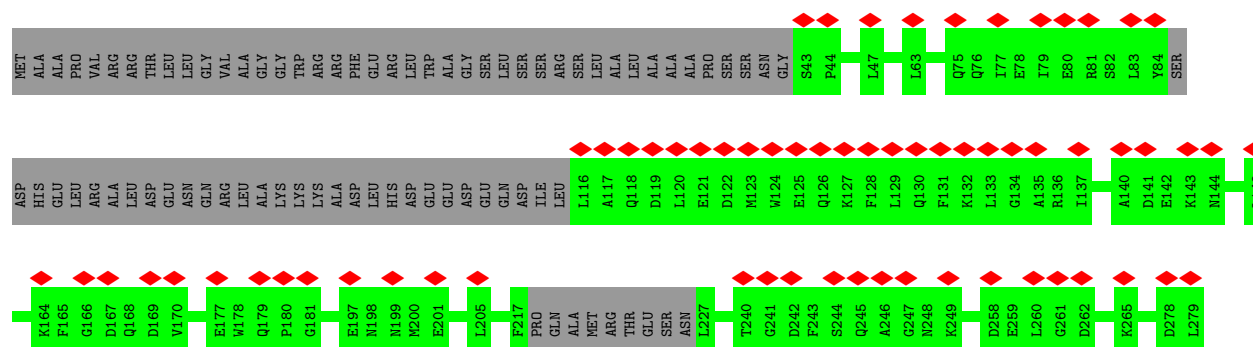
- Molecule 40: 39S ribosomal protein L45, mitochondrial

Chain d:  64% 35%



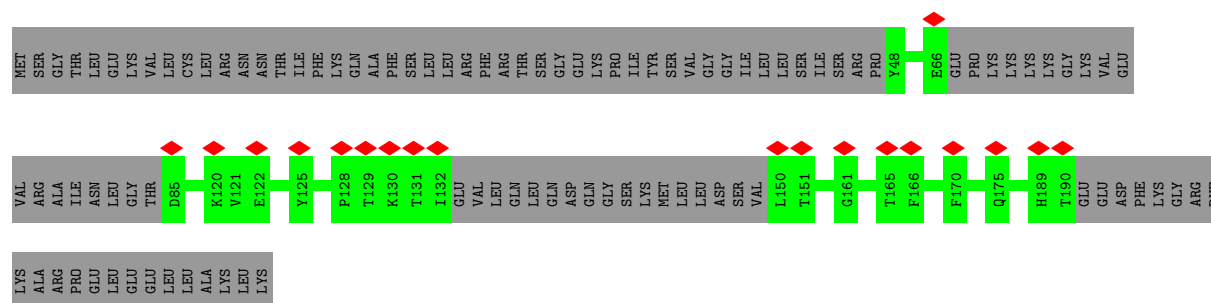
- Molecule 41: 39S ribosomal protein L46, mitochondrial

Chain e: 23% 71% 29%



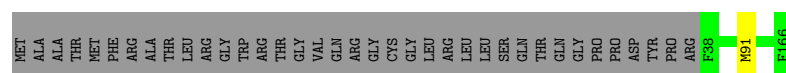
- Molecule 42: 39S ribosomal protein L48, mitochondrial

Chain f: 9% 51% 49%



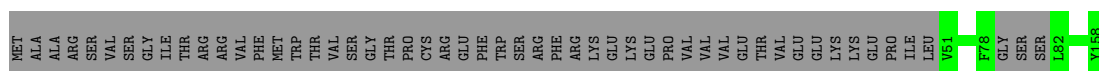
- Molecule 43: 39S ribosomal protein L49, mitochondrial

Chain g: 77% 22%



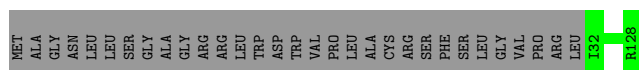
- Molecule 44: 39S ribosomal protein L50, mitochondrial

Chain h: 66% 34%



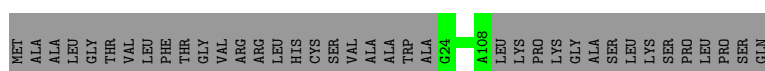
- Molecule 45: 39S ribosomal protein L51, mitochondrial

Chain i: 76% 24%



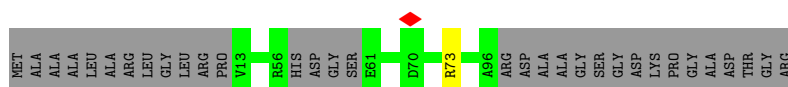
- Molecule 46: 39S ribosomal protein L52, mitochondrial

Chain j: 69% 31%



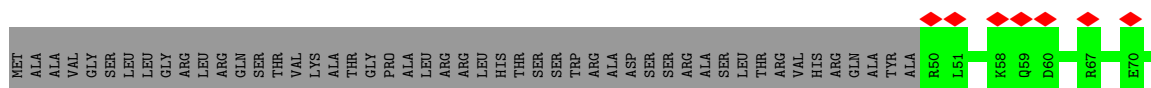
- Molecule 47: 39S ribosomal protein L53, mitochondrial

Chain k: 71% 29%



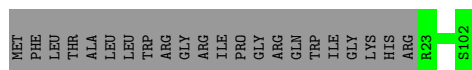
- Molecule 48: 39S ribosomal protein L55, mitochondrial

Chain m: 6% 23% 77%



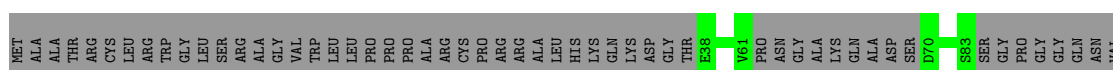
- Molecule 49: Ribosomal protein 63, mitochondrial

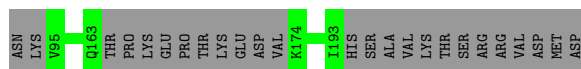
Chain o: 78% 22%



- Molecule 50: Peptidyl-tRNA hydrolase ICT1, mitochondrial

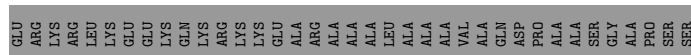
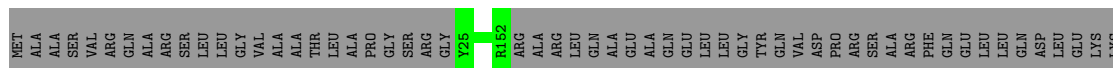
Chain p: 62% 38%





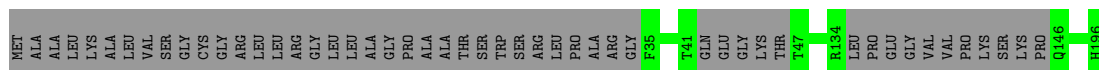
- Molecule 51: Growth arrest and DNA damage-inducible proteins-interacting protein 1

Chain q: 58% 42%



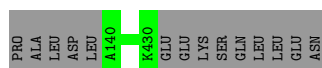
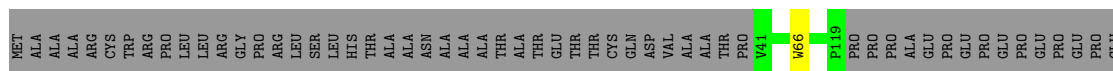
- Molecule 52: 39S ribosomal protein S18a, mitochondrial

Chain r: 74% 26%



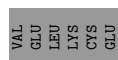
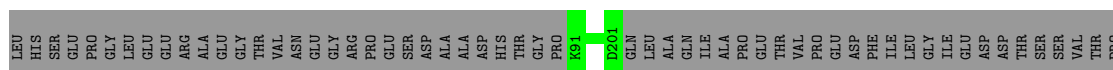
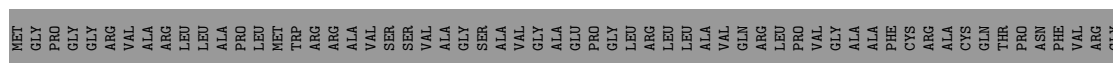
- Molecule 53: 39S ribosomal protein S30, mitochondrial

Chain s: 84% 16%



- Molecule 54: Mitochondrial assembly of ribosomal large subunit protein 1

Chain u: 47% 53%



- Molecule 55: MIEF1 upstream open reading frame protein

Chain v: 96% ..



- Molecule 56: Acyl carrier protein, mitochondrial





MET	ALA	SER	ARG	VAL	LEU	SER	ALA	TYR	VAL	SER	ARG	LEU	PRO	ALA	ALA	PHE	ALA	PRO	LEU	PRO	ARG	VAL	ARG	ARG	MET	LEU	ALA	VAL	ALA	ALA	ARG	PRO	LEU	SER	THR	ALA	ALA	LEU	CYS	SER	ALA	GLY	THR	GLN	THR	ARG	LEU	GLY	THR	LEU	LEU	GLN	PRO	ALA	VAL	LEU	ALA	GLN	VAL	PRO	GLY	ARG
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VAL	THR	GLN	LEU	CYS	ARG	GLN	TYR	SER	ASP	MET	PRO	PRO	L74	K106	K152	ASP	VAL	TYR	GLU
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## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	1060638	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 ( $e^-/\text{\AA}^2$ )	36	Depositor
Minimum defocus (nm)	300	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.128	Depositor
Minimum map value	-0.035	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.004	Depositor
Map size (Å)	367.49997, 367.49997, 367.49997	wwPDB
Map dimensions	350, 350, 350	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, ZN

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	0.33	0/895	0.53	0/1201
2	1	0.24	0/438	0.54	0/583
3	2	0.37	0/373	0.57	0/496
4	3	0.39	0/852	0.53	0/1136
5	4	0.36	0/341	0.57	0/451
6	5	0.32	0/3294	0.50	0/4488
7	6	0.31	0/2809	0.53	0/3818
8	7	0.30	0/2391	0.47	0/3234
9	8	0.25	0/665	0.49	0/894
10	9	0.34	0/972	0.49	0/1306
11	A	0.65	0/32679	0.83	23/50833 (0.0%)
12	B	0.22	0/1328	0.76	0/2056
13	C	0.25	0/2655	0.49	0/3601
14	D	0.34	0/1910	0.57	0/2569
15	E	0.38	0/2497	0.49	0/3386
16	F	0.37	0/2071	0.52	0/2817
17	G	0.27	0/1974	0.47	0/2652
18	H	0.31	0/798	0.53	0/1073
19	I	0.26	0/1252	0.50	0/1686
20	J	0.24	0/1077	0.50	0/1452
21	K	0.38	0/1495	0.49	0/2029
22	L	0.33	0/904	0.55	0/1218
23	M	0.37	0/2359	0.53	0/3185
24	N	0.28	0/1533	0.52	0/2061
25	O	0.34	0/1269	0.54	0/1708
26	P	0.30	0/1173	0.53	0/1588
27	Q	0.34	0/1846	0.51	0/2487
28	R	0.40	0/987	0.56	0/1320
29	S	0.36	0/1276	0.54	0/1729
30	T	0.40	0/1402	0.51	0/1886
31	U	0.37	0/1183	0.55	0/1600
32	V	0.33	0/404	0.44	0/545

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
33	W	0.38	0/823	0.50	0/1113
34	X	0.34	0/2090	0.47	0/2825
35	Y	0.34	0/1552	0.50	0/2079
36	Z	0.32	0/960	0.49	0/1295
37	a	0.36	0/616	0.50	0/833
38	b	0.36	0/1202	0.57	0/1626
39	c	0.33	0/2264	0.48	0/3059
40	d	0.26	0/1702	0.47	0/2307
41	e	0.24	0/1633	0.50	0/2204
42	f	0.27	0/873	0.44	0/1180
43	g	0.38	0/1102	0.51	0/1503
44	h	0.27	0/884	0.47	0/1203
45	i	0.39	0/849	0.54	0/1135
46	j	0.31	0/698	0.50	0/940
47	k	0.25	0/635	0.54	0/855
48	m	0.23	0/250	0.64	0/336
49	o	0.33	0/693	0.52	0/931
50	p	0.27	0/1071	0.52	0/1433
51	q	0.28	0/1107	0.50	0/1498
52	r	0.33	0/1238	0.52	0/1676
53	s	0.37	0/3114	0.52	0/4225
54	u	0.29	0/949	0.51	0/1281
55	v	0.26	0/597	0.60	0/796
56	w	0.25	0/647	0.46	0/871
All	All	0.45	0/104651	0.64	23/148292 (0.0%)

There are no bond length outliers.

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	A	2523	C	C2-N1-C1'	8.72	128.40	118.80
11	A	2523	C	N1-C2-O2	7.99	123.70	118.90
11	A	2523	C	N3-C2-O2	-6.47	117.37	121.90
11	A	2543	C	C2'-C3'-O3'	6.43	123.99	113.70
11	A	2523	C	C6-N1-C1'	-6.22	113.33	120.80

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	0	106/188 (56%)	104 (98%)	2 (2%)	0	100	100
2	1	50/65 (77%)	48 (96%)	2 (4%)	0	100	100
3	2	43/92 (47%)	42 (98%)	1 (2%)	0	100	100
4	3	93/188 (50%)	91 (98%)	2 (2%)	0	100	100
5	4	35/103 (34%)	34 (97%)	1 (3%)	0	100	100
6	5	390/423 (92%)	382 (98%)	8 (2%)	0	100	100
7	6	316/380 (83%)	305 (96%)	11 (4%)	0	100	100
8	7	285/338 (84%)	271 (95%)	14 (5%)	0	100	100
9	8	75/206 (36%)	73 (97%)	2 (3%)	0	100	100
10	9	113/137 (82%)	111 (98%)	2 (2%)	0	100	100
13	C	324/384 (84%)	312 (96%)	12 (4%)	0	100	100
14	D	238/305 (78%)	230 (97%)	8 (3%)	0	100	100
15	E	306/348 (88%)	300 (98%)	6 (2%)	0	100	100
16	F	248/311 (80%)	245 (99%)	3 (1%)	0	100	100
17	G	236/381 (62%)	227 (96%)	9 (4%)	0	100	100
18	H	93/267 (35%)	88 (95%)	5 (5%)	0	100	100
19	I	147/261 (56%)	140 (95%)	7 (5%)	0	100	100
20	J	138/192 (72%)	132 (96%)	6 (4%)	0	100	100
21	K	175/178 (98%)	172 (98%)	3 (2%)	0	100	100
22	L	113/145 (78%)	112 (99%)	1 (1%)	0	100	100
23	M	285/296 (96%)	282 (99%)	3 (1%)	0	100	100

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
24	N	177/251 (70%)	174 (98%)	3 (2%)	0	100	100
25	O	150/175 (86%)	146 (97%)	4 (3%)	0	100	100
26	P	139/180 (77%)	132 (95%)	7 (5%)	0	100	100
27	Q	215/292 (74%)	211 (98%)	4 (2%)	0	100	100
28	R	114/149 (76%)	113 (99%)	1 (1%)	0	100	100
29	S	154/205 (75%)	151 (98%)	3 (2%)	0	100	100
30	T	164/206 (80%)	162 (99%)	2 (1%)	0	100	100
31	U	135/153 (88%)	134 (99%)	1 (1%)	0	100	100
32	V	45/216 (21%)	44 (98%)	1 (2%)	0	100	100
33	W	98/148 (66%)	98 (100%)	0	0	100	100
34	X	241/256 (94%)	238 (99%)	3 (1%)	0	100	100
35	Y	174/250 (70%)	171 (98%)	3 (2%)	0	100	100
36	Z	113/161 (70%)	111 (98%)	2 (2%)	0	100	100
37	a	67/142 (47%)	67 (100%)	0	0	100	100
38	b	146/215 (68%)	141 (97%)	5 (3%)	0	100	100
39	c	271/332 (82%)	268 (99%)	3 (1%)	0	100	100
40	d	189/306 (62%)	186 (98%)	3 (2%)	0	100	100
41	e	191/279 (68%)	175 (92%)	16 (8%)	0	100	100
42	f	102/212 (48%)	92 (90%)	10 (10%)	0	100	100
43	g	127/166 (76%)	121 (95%)	6 (5%)	0	100	100
44	h	101/158 (64%)	98 (97%)	3 (3%)	0	100	100
45	i	95/128 (74%)	94 (99%)	1 (1%)	0	100	100
46	j	83/123 (68%)	82 (99%)	1 (1%)	0	100	100
47	k	76/112 (68%)	71 (93%)	5 (7%)	0	100	100
48	m	27/128 (21%)	19 (70%)	8 (30%)	0	100	100
49	o	78/102 (76%)	78 (100%)	0	0	100	100
50	p	119/206 (58%)	116 (98%)	3 (2%)	0	100	100
51	q	126/222 (57%)	126 (100%)	0	0	100	100
52	r	140/196 (71%)	136 (97%)	4 (3%)	0	100	100
53	s	366/439 (83%)	363 (99%)	3 (1%)	0	100	100
54	u	109/234 (47%)	103 (94%)	6 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
55	v	67/70 (96%)	66 (98%)	1 (2%)	0	100	100
56	w	77/156 (49%)	70 (91%)	7 (9%)	0	100	100
All	All	8285/11756 (70%)	8058 (97%)	227 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	49/60 (82%)	49 (100%)	0	100	100
3	2	39/72 (54%)	39 (100%)	0	100	100
4	3	88/166 (53%)	88 (100%)	0	100	100
5	4	36/89 (40%)	36 (100%)	0	100	100
6	5	353/368 (96%)	351 (99%)	2 (1%)	84	91
7	6	286/332 (86%)	285 (100%)	1 (0%)	91	96
8	7	263/303 (87%)	258 (98%)	5 (2%)	52	67
9	8	70/190 (37%)	70 (100%)	0	100	100
10	9	99/112 (88%)	99 (100%)	0	100	100
13	C	284/328 (87%)	284 (100%)	0	100	100
14	D	194/245 (79%)	192 (99%)	2 (1%)	73	84
15	E	262/290 (90%)	262 (100%)	0	100	100
16	F	217/262 (83%)	216 (100%)	1 (0%)	86	93
17	G	221/350 (63%)	221 (100%)	0	100	100
18	H	86/228 (38%)	85 (99%)	1 (1%)	67	80
19	I	139/232 (60%)	139 (100%)	0	100	100
20	J	113/150 (75%)	113 (100%)	0	100	100
21	K	155/156 (99%)	155 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
22	L	98/124 (79%)	97 (99%)	1 (1%)	73	84
23	M	245/249 (98%)	244 (100%)	1 (0%)	89	95
24	N	156/211 (74%)	156 (100%)	0	100	100
25	O	133/150 (89%)	133 (100%)	0	100	100
26	P	123/155 (79%)	123 (100%)	0	100	100
27	Q	199/256 (78%)	199 (100%)	0	100	100
28	R	101/126 (80%)	101 (100%)	0	100	100
29	S	141/180 (78%)	141 (100%)	0	100	100
30	T	146/176 (83%)	146 (100%)	0	100	100
31	U	124/135 (92%)	124 (100%)	0	100	100
32	V	43/191 (22%)	43 (100%)	0	100	100
33	W	83/119 (70%)	83 (100%)	0	100	100
34	X	219/229 (96%)	218 (100%)	1 (0%)	86	93
35	Y	159/223 (71%)	159 (100%)	0	100	100
36	Z	106/147 (72%)	105 (99%)	1 (1%)	75	86
37	a	67/133 (50%)	67 (100%)	0	100	100
38	b	130/186 (70%)	130 (100%)	0	100	100
39	c	241/288 (84%)	241 (100%)	0	100	100
40	d	184/274 (67%)	182 (99%)	2 (1%)	70	82
41	e	171/236 (72%)	171 (100%)	0	100	100
42	f	95/188 (50%)	95 (100%)	0	100	100
43	g	119/148 (80%)	118 (99%)	1 (1%)	79	88
44	h	100/148 (68%)	100 (100%)	0	100	100
45	i	86/110 (78%)	86 (100%)	0	100	100
46	j	68/97 (70%)	68 (100%)	0	100	100
47	k	71/90 (79%)	70 (99%)	1 (1%)	62	77
48	m	27/113 (24%)	27 (100%)	0	100	100
49	o	69/87 (79%)	69 (100%)	0	100	100
50	p	117/181 (65%)	117 (100%)	0	100	100
51	q	110/178 (62%)	110 (100%)	0	100	100
52	r	133/169 (79%)	133 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
53	s	326/381 (86%)	325 (100%)	1 (0%)	91	96
54	u	105/200 (52%)	105 (100%)	0	100	100
55	v	59/60 (98%)	57 (97%)	2 (3%)	32	42
56	w	73/136 (54%)	72 (99%)	1 (1%)	62	77
All	All	7478/10171 (74%)	7454 (100%)	24 (0%)	90	96

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
34	X	23	ARG
40	d	198	ARG
40	d	51	LYS
43	g	91	MET
8	7	234	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
41	e	212	HIS
38	b	58	ASN
16	F	241	ASN
14	D	271	ASN
29	S	118	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
11	A	1357/1559 (87%)	323 (23%)	17 (1%)
12	B	51/69 (73%)	16 (31%)	1 (1%)
All	All	1408/1628 (86%)	339 (24%)	18 (1%)

5 of 339 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
11	A	1677	C
11	A	1678	C
11	A	1679	U
11	A	1680	A
11	A	1681	G

5 of 18 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
11	A	3072	U
12	B	1607	U
11	A	3196	G
11	A	2530	A
11	A	3041	U

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 64 ligands modelled in this entry, 64 are monoatomic - leaving 0 for Mogul analysis.

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.

## 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.

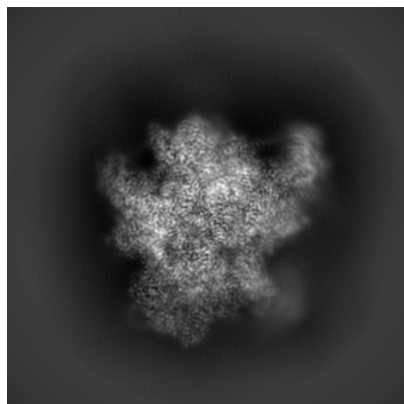
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12865. These allow visual inspection of the internal detail of the map and identification of artifacts.

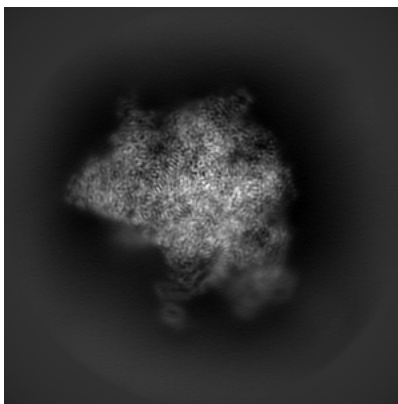
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

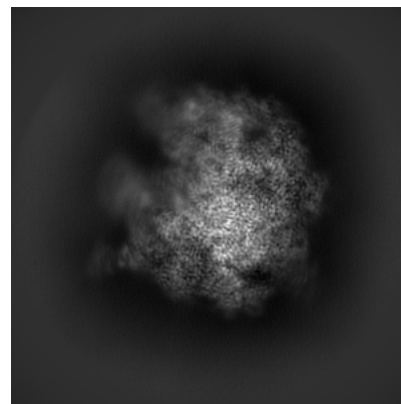
#### 6.1.1 Primary map



X

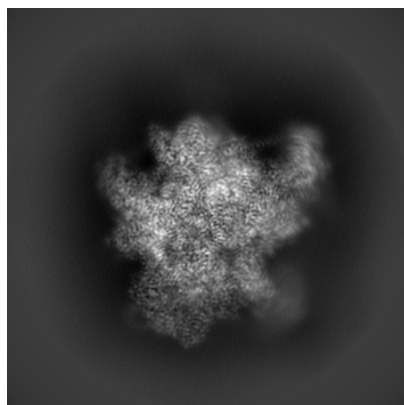


Y

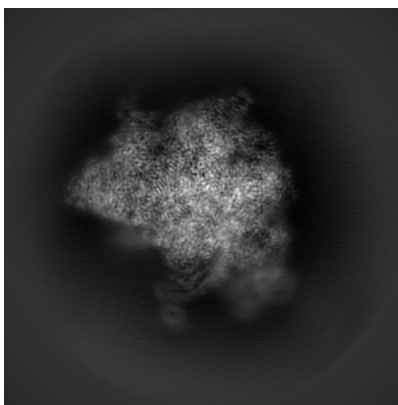


Z

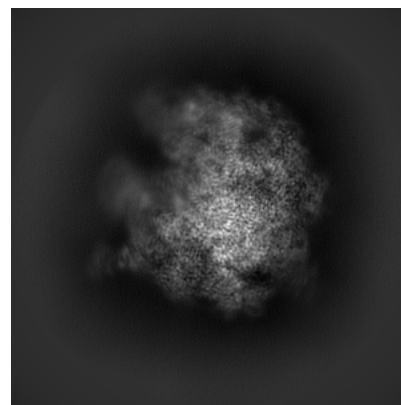
#### 6.1.2 Raw map



X



Y

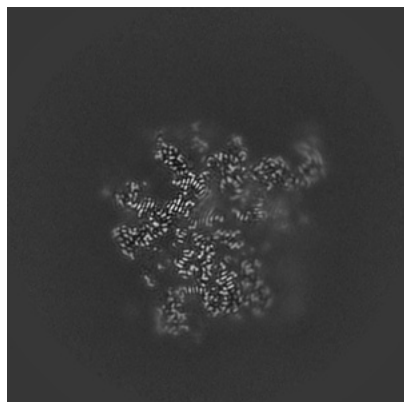


Z

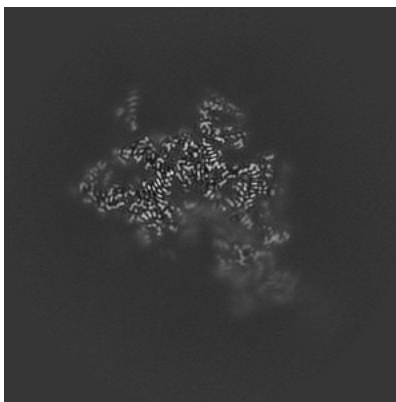
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

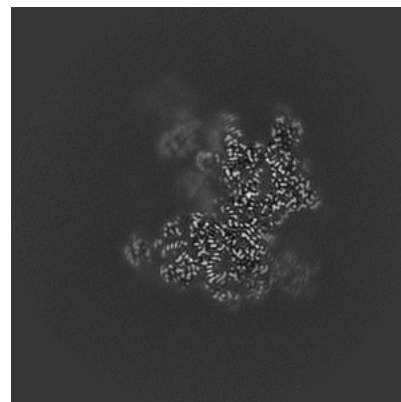
### 6.2.1 Primary map



X Index: 175

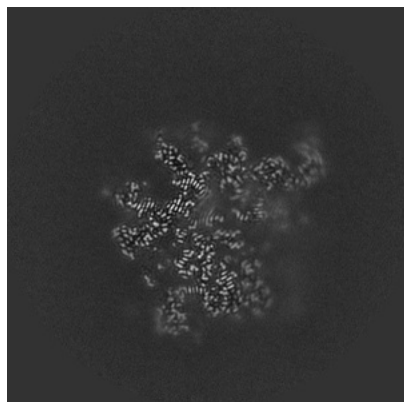


Y Index: 175

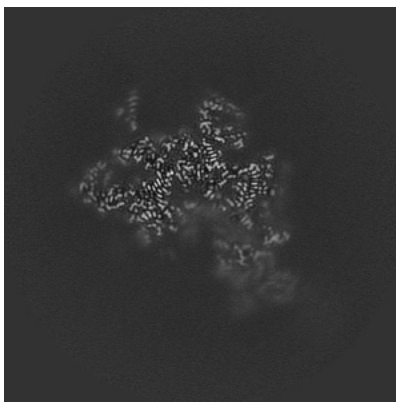


Z Index: 175

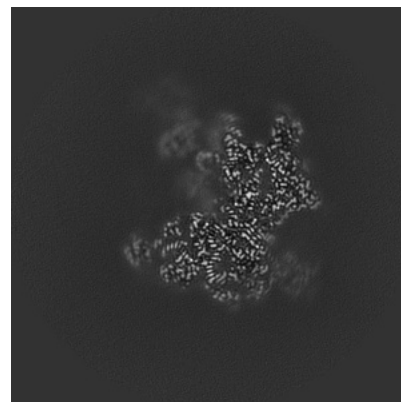
### 6.2.2 Raw map



X Index: 175



Y Index: 175

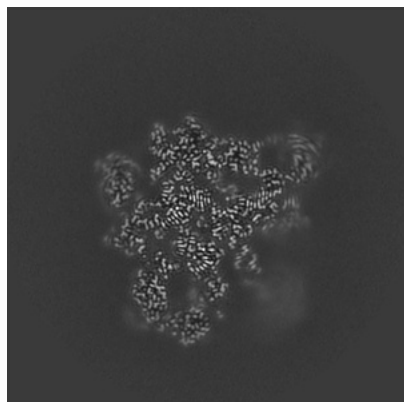


Z Index: 175

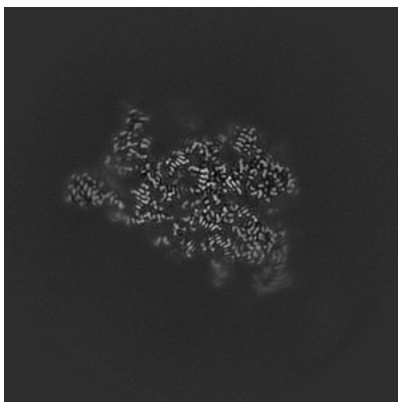
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

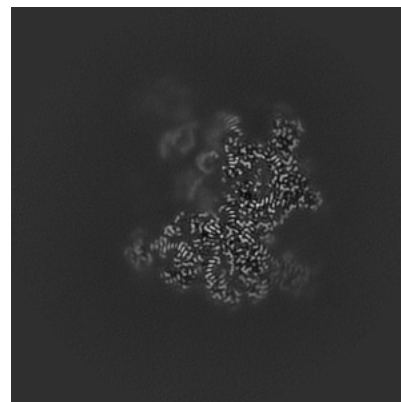
### 6.3.1 Primary map



X Index: 192

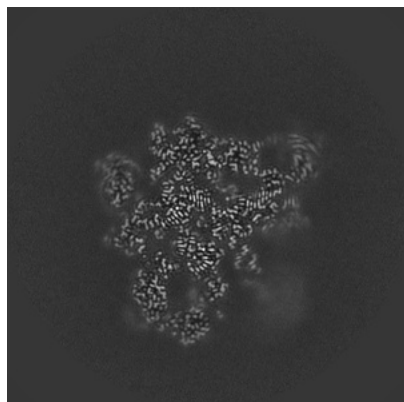


Y Index: 159

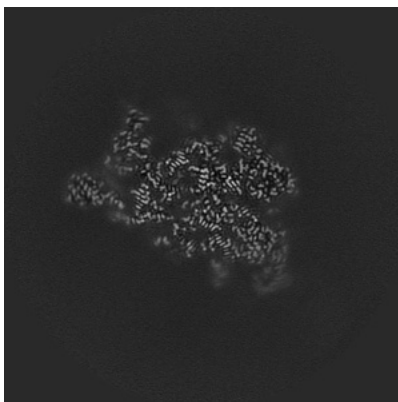


Z Index: 177

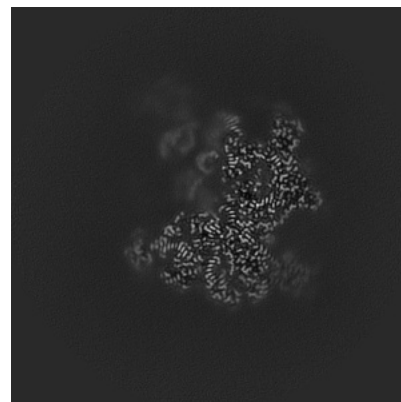
### 6.3.2 Raw map



X Index: 192



Y Index: 159



Z Index: 177

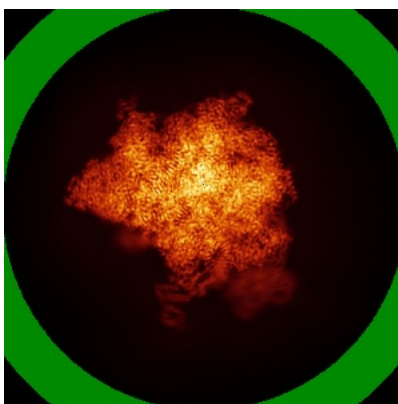
The images above show the largest variance slices of the map in three orthogonal directions.

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

### 6.4.1 Primary map



X

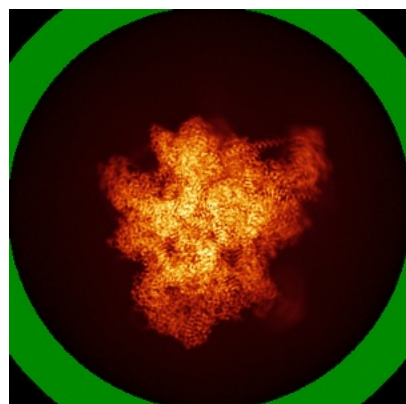


Y

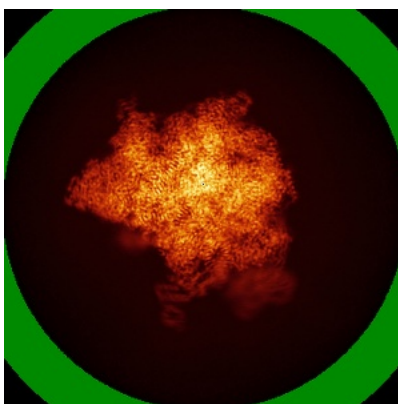


Z

### 6.4.2 Raw map



X



Y



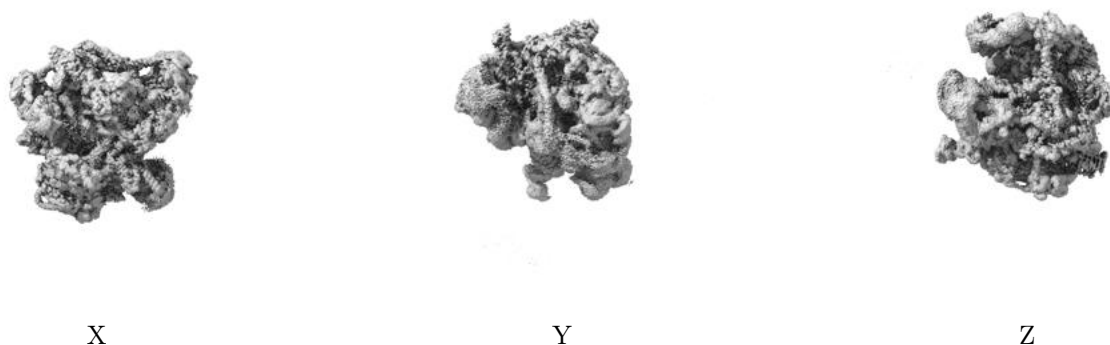
Z

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



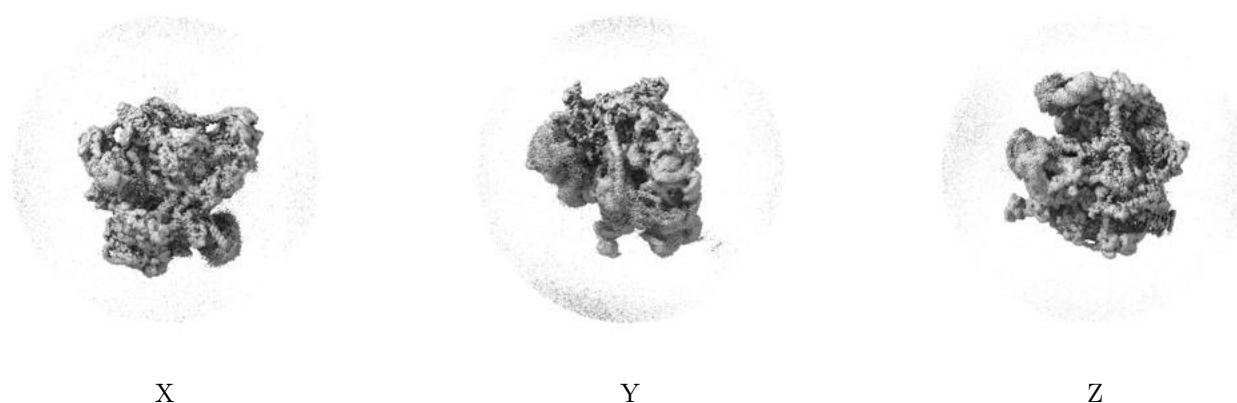
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.004. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



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

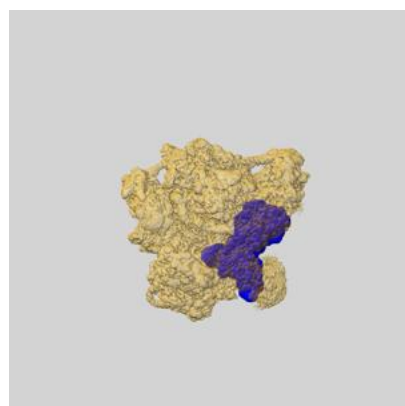
## 6.6 Mask visualisation [i](#)

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

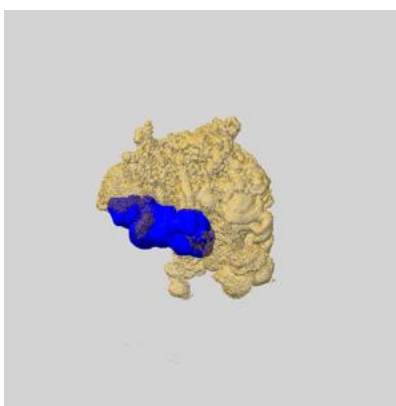
A mask typically either:

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

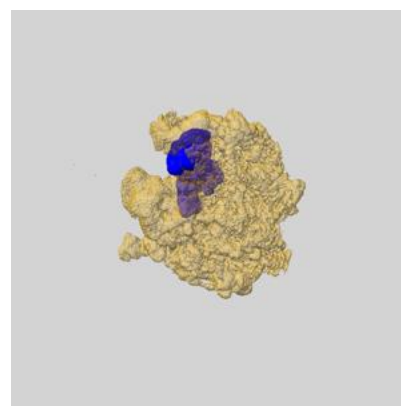
### 6.6.1 emd\_12865\_msk\_1.map [i](#)



X

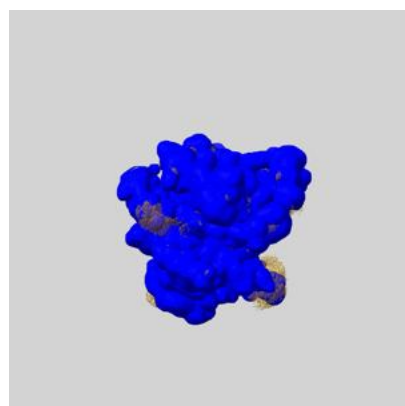


Y

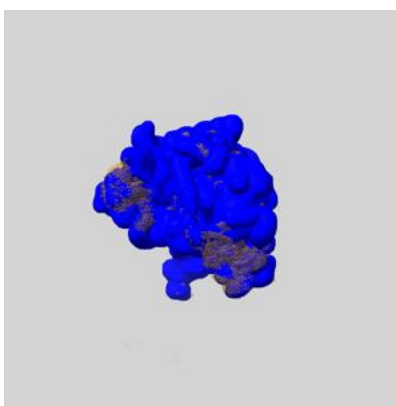


Z

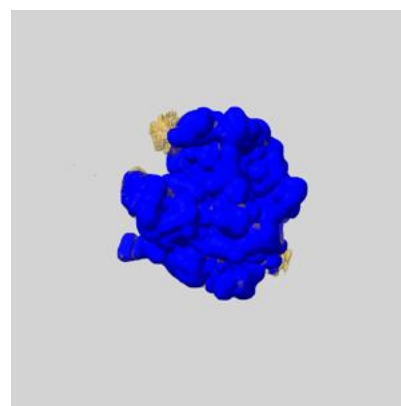
### 6.6.2 emd\_12865\_msk\_2.map [i](#)



X



Y



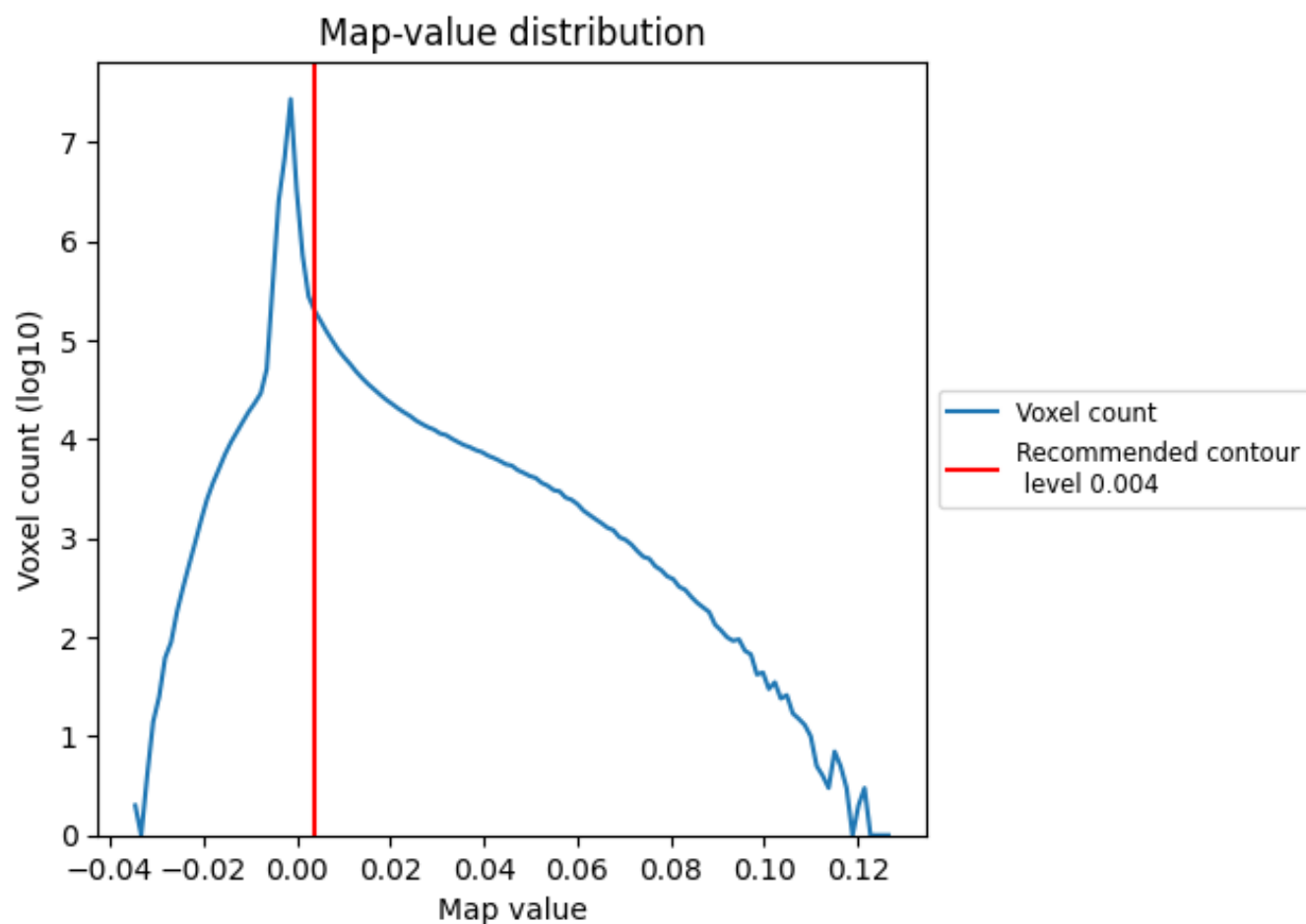
Z



## 7 Map analysis [i](#)

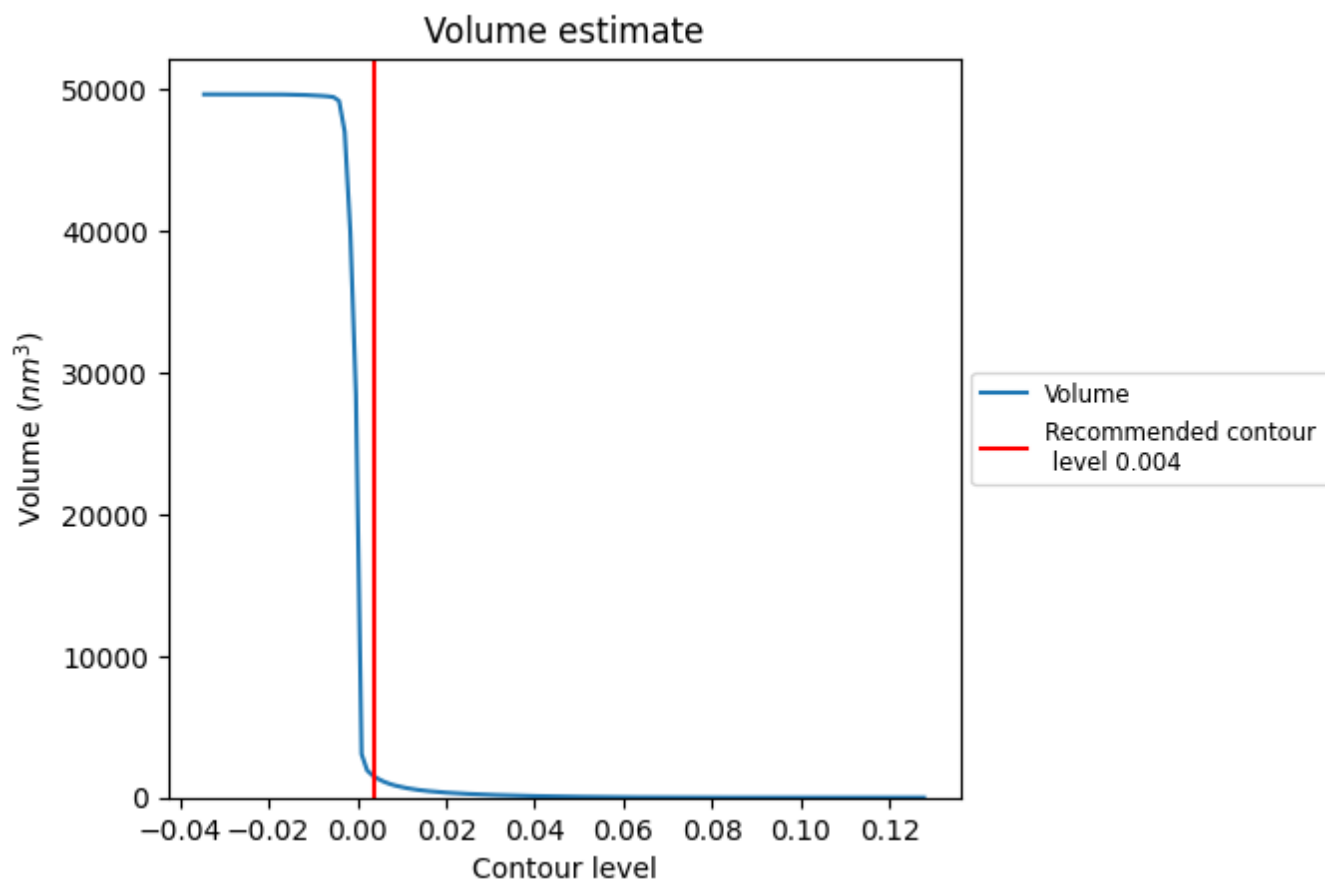
This section contains the results of statistical analysis of the map.

### 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.

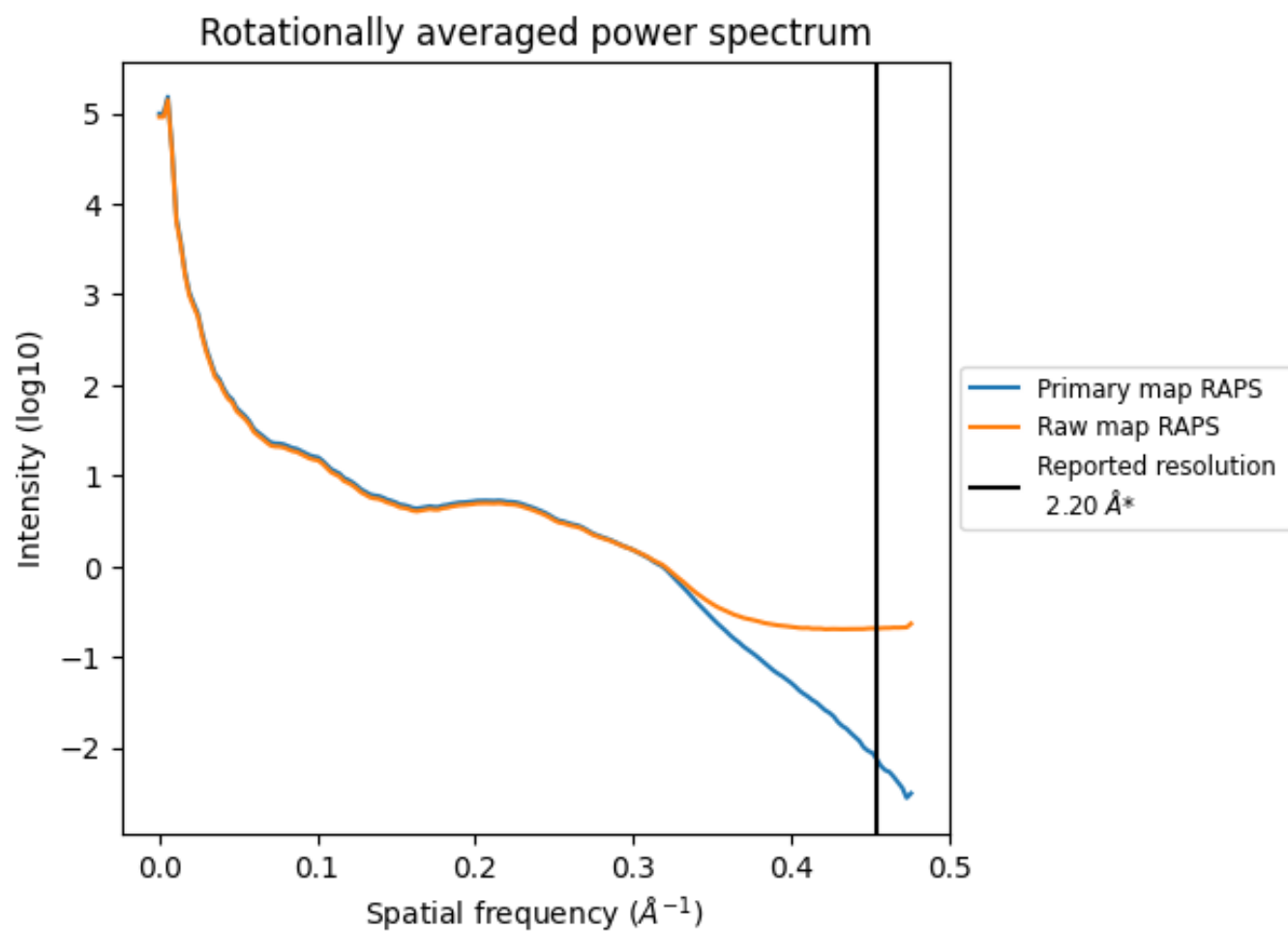
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1458 nm<sup>3</sup>; this corresponds to an approximate mass of 1317 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 ⓘ

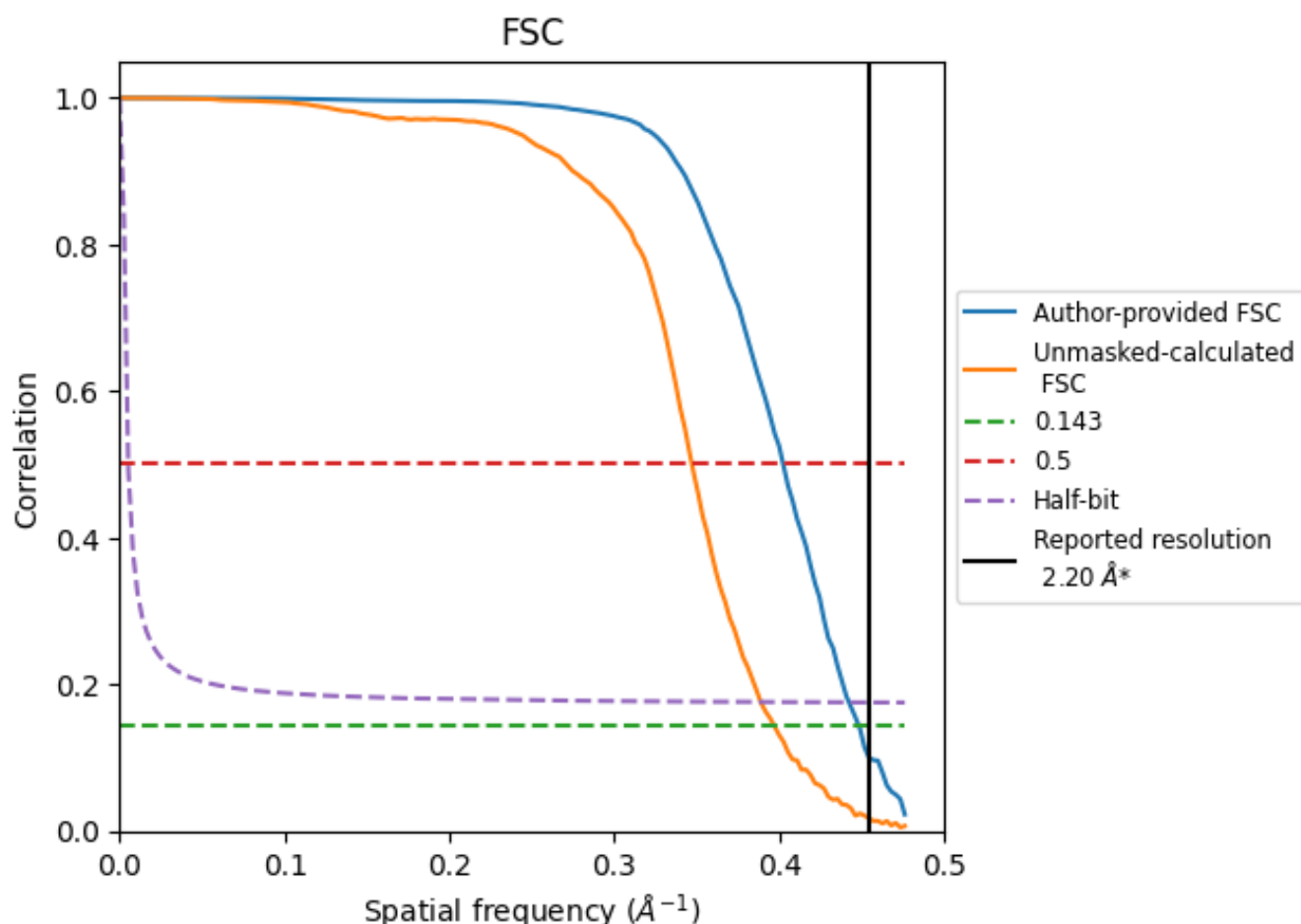


\*Reported resolution corresponds to spatial frequency of 0.455  $\text{\AA}^{-1}$

## 8 Fourier-Shell correlation [i](#)

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

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.455 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

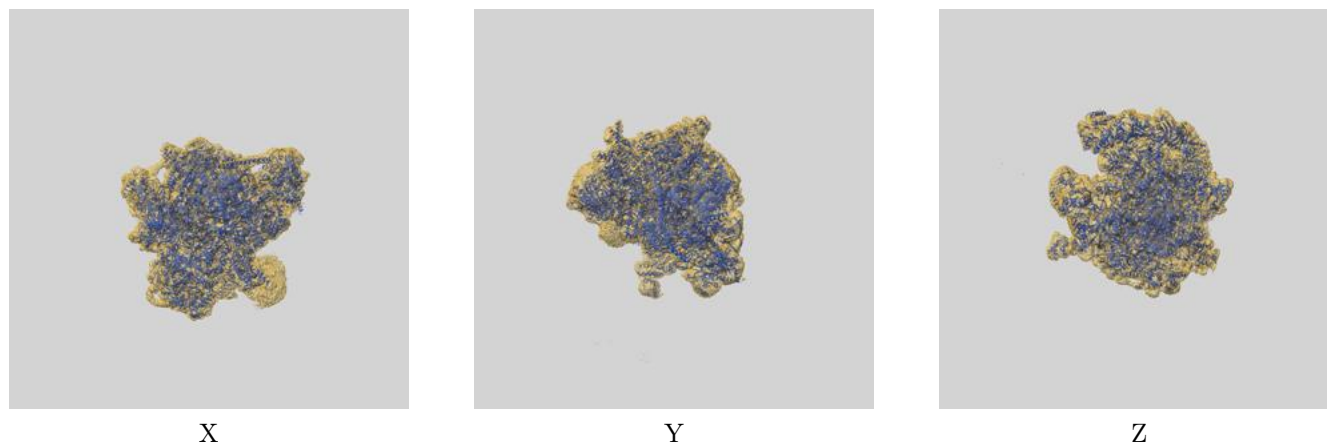
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.20	-	-
Author-provided FSC curve	2.23	2.48	2.26
Unmasked-calculated*	2.52	2.88	2.57

\*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 2.52 differs from the reported value 2.2 by more than 10 %

## 9 Map-model fit [i](#)

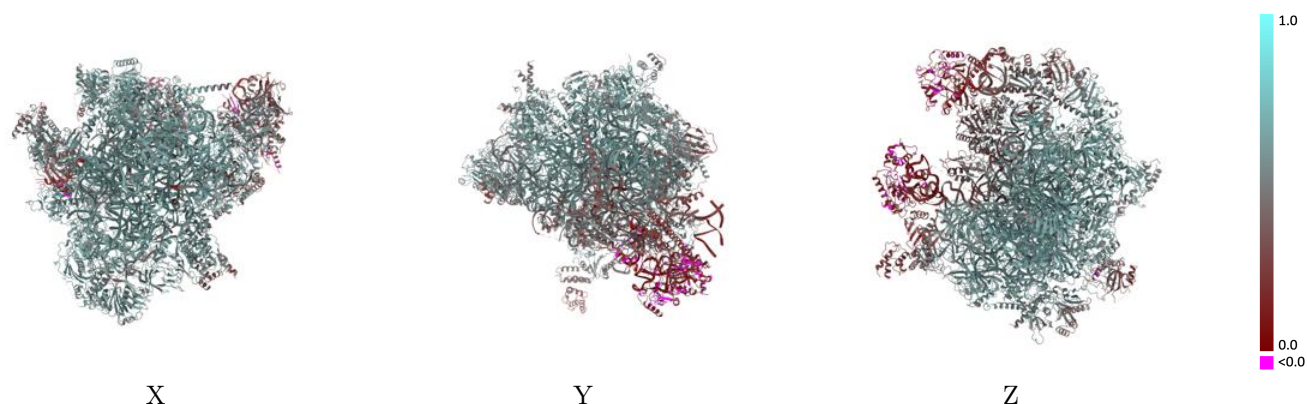
This section contains information regarding the fit between EMDB map EMD-12865 and PDB model 7OF0. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

### 9.1 Map-model overlay [i](#)



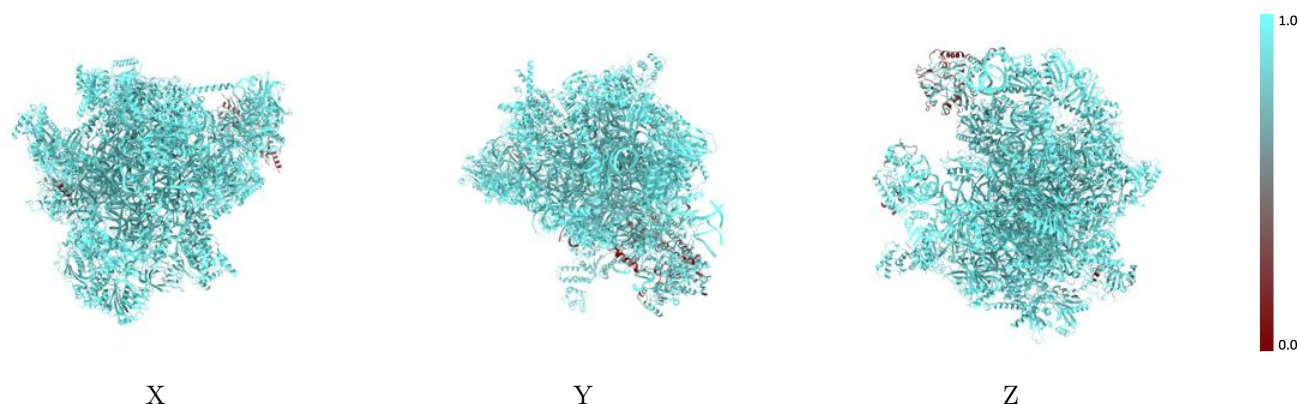
The images above show the 3D surface view of the map at the recommended contour level 0.004 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



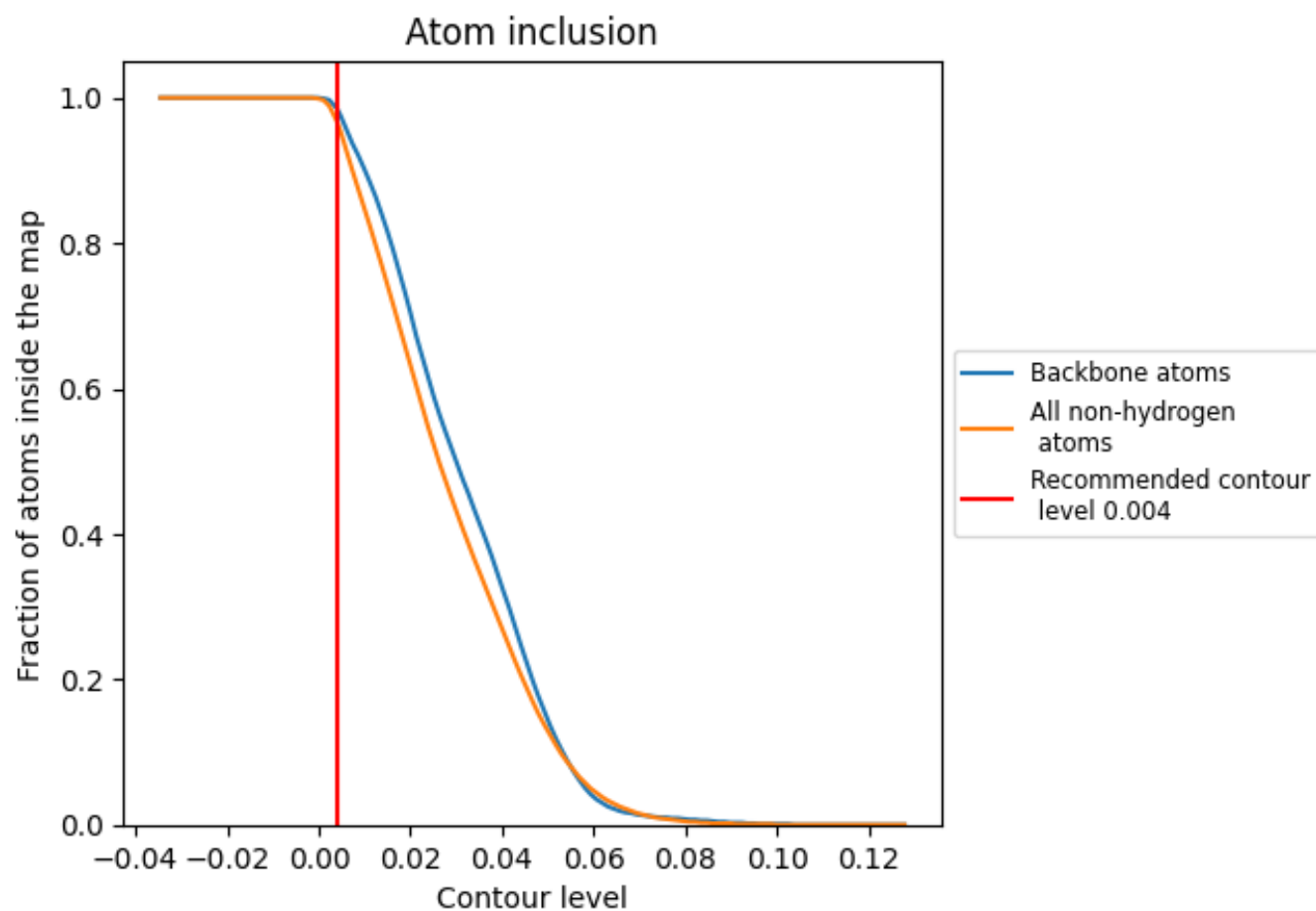
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.004).

## 9.4 Atom inclusion [i](#)









































































At the recommended contour level, 98% of all backbone atoms, 97% 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 (0.004) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9670	 0.5360
0	 0.9720	 0.5920
1	 0.9670	 0.3880
2	 1.0000	 0.6690
3	 0.9930	 0.6490
4	 1.0000	 0.5770
5	 0.9850	 0.5650
6	 0.9780	 0.4590
7	 0.9800	 0.5180
8	 0.5940	 0.1780
9	 0.9810	 0.5830
A	 0.9950	 0.5770
B	 0.9760	 0.2960
C	 0.9710	 0.3940
D	 0.9800	 0.6020
E	 0.9900	 0.6120
F	 0.9950	 0.6350
G	 0.9540	 0.4480
H	 0.9720	 0.5520
I	 0.9400	 0.2380
J	 0.8390	 0.0850
K	 0.9940	 0.6290
L	 0.9890	 0.5960
M	 0.9900	 0.6230
N	 0.9700	 0.4400
O	 0.9910	 0.6230
P	 0.9880	 0.5190
Q	 0.9840	 0.5940
R	 0.9900	 0.6430
S	 0.9870	 0.6140
T	 0.9850	 0.6350
U	 0.8900	 0.5600
V	 0.9790	 0.5590
W	 0.9910	 0.6110
X	 0.9800	 0.5940



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Chain	Atom inclusion	Q-score
Y	 0.9820	 0.6120
Z	 0.9890	 0.6050
a	 0.9880	 0.6180
b	 0.9900	 0.6230
c	 0.9800	 0.5840
d	 0.8620	 0.3900
e	 0.5420	 0.1240
f	 0.6860	 0.3110
g	 0.9920	 0.6180
h	 0.9800	 0.5220
i	 0.9980	 0.6570
j	 0.9890	 0.5920
k	 0.9590	 0.2990
m	 0.5790	 0.1720
o	 0.9880	 0.6200
p	 0.9630	 0.5160
q	 0.9540	 0.4880
r	 0.9910	 0.5620
s	 0.9870	 0.6100
u	 0.9740	 0.4640
v	 0.9540	 0.3520
w	 0.9100	 0.2280