



# wwPDB EM Validation Summary Report ⓘ

Mar 19, 2025 – 06:50 PM EDT

PDB ID : 9B1X  
EMDB ID : EMD-44091  
Title : HWS19 strain gidB mutant mycobacterial ribosome  
Authors : Chen, Y.; Young, I.D.; Fraser, J.S.; Javid, B.  
Deposited on : 2024-03-14  
Resolution : 3.07 Å(reported)  
Based on initial models : 5o60, 5o5j

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.dev117  
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) : 2.41.4

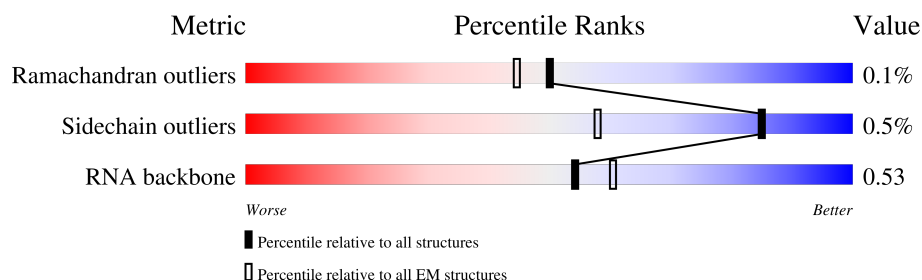
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.07 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.




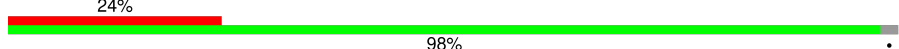
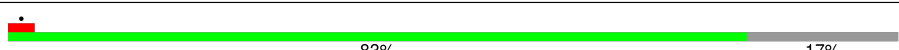
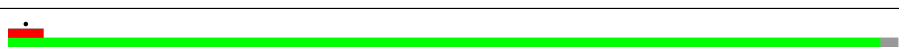
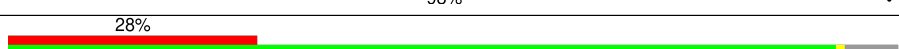
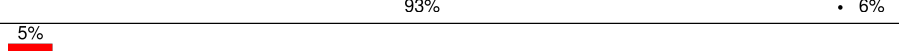
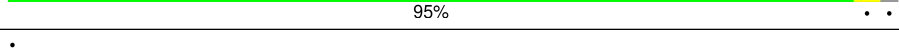
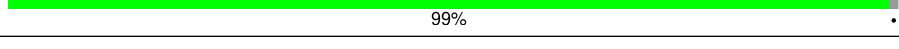

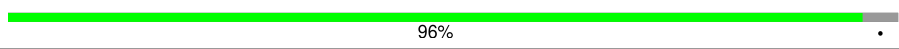


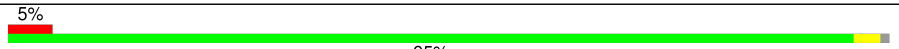


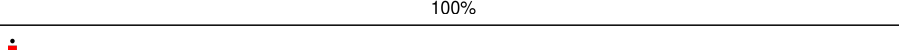


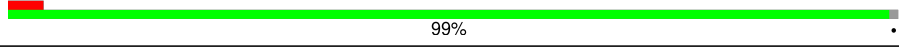
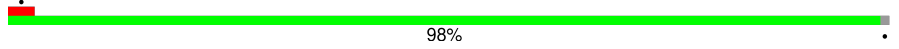
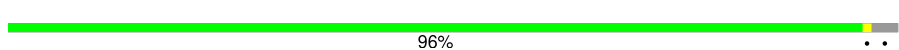
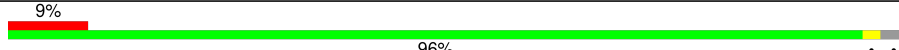

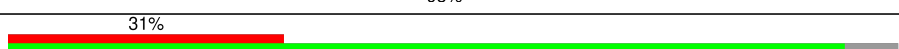
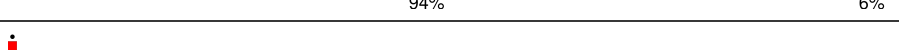
Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

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	A	1528	
2	B	33	
3	C	275	
4	D	201	
5	E	214	
6	F	96	
7	G	156	
8	H	132	

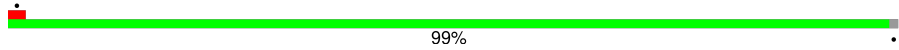

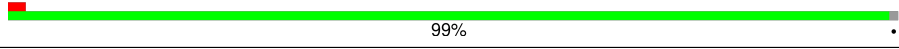
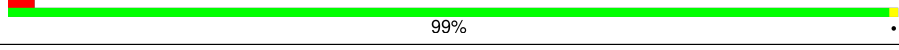
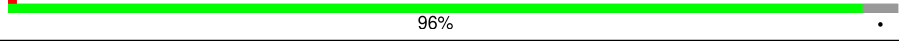
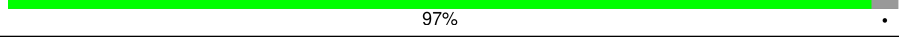

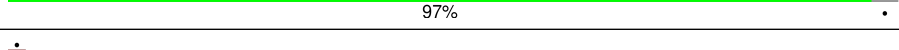
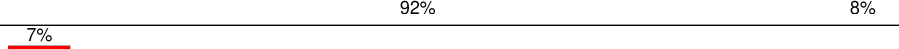
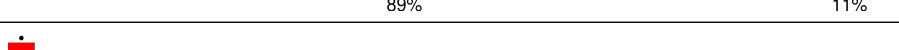

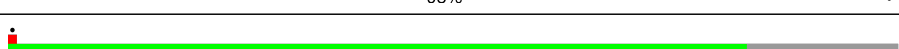


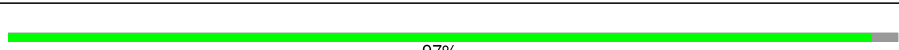
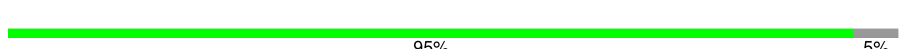

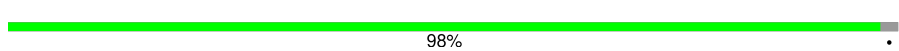
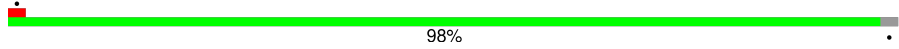
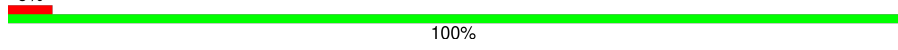

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Mol	Chain	Length	Quality of chain
9	I	150	
10	J	101	
11	K	138	
12	L	124	
13	M	124	
14	N	61	
15	O	89	
16	P	156	
17	Q	98	
18	R	84	
19	S	93	
20	T	86	
21	V	277	
22	X	6	
23	Y	3120	
24	U	118	
25	Z	278	
26	a	217	
27	b	215	
28	c	187	
29	d	179	
30	e	142	
31	f	147	
32	g	122	
33	h	147	

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Mol	Chain	Length	Quality of chain
34	i	138	
35	j	199	
36	k	127	
37	l	113	
38	m	129	
39	n	103	
40	o	153	
41	p	100	
42	q	105	
43	r	215	
44	s	88	
45	t	64	
46	u	77	
47	5	24	
48	v	175	
49	w	61	
50	x	57	
51	y	55	
52	z	47	
53	1	64	
54	2	37	

## 2 Entry composition

There are 56 unique types of molecules in this entry. The entry contains 242013 atoms, of which 97195 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 16S rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	1511	Total	C	H	N	O	P	0	0
			48755	14448	16316	5930	10550	1511		

- Molecule 2 is a protein called Conserved domain protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	32	Total	C	H	N	O	S	0	0
			623	172	343	71	36	1		

- Molecule 3 is a protein called Small ribosomal subunit protein uS3.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	208	Total	C	H	N	O	S	0	0
			3368	1036	1708	322	298	4		

- Molecule 4 is a protein called Small ribosomal subunit protein uS4.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	200	Total	C	H	N	O	S	0	0
			3310	1028	1669	316	295	2		

- Molecule 5 is a protein called Small ribosomal subunit protein uS5.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	E	180	Total	C	H	N	O	S	0	0
			2657	812	1361	245	235	4		

- Molecule 6 is a protein called Small ribosomal subunit protein bS6.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	F	96	Total	C	H	N	O	S	0	0
			1568	486	797	138	145	2		

- Molecule 7 is a protein called Small ribosomal subunit protein uS7.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	G	155	Total	C	H	N	O	S	0	0
			2514	768	1282	241	221	2		

- Molecule 8 is a protein called Small ribosomal subunit protein uS8.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	H	131	Total	C	H	N	O	S	0	0
			2057	633	1047	189	187	1		

- Molecule 9 is a protein called Small ribosomal subunit protein uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	126	Total	C	H	N	O	0	0
			2044	630	1050	194	170		

- Molecule 10 is a protein called Small ribosomal subunit protein uS10.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	J	99	Total	C	H	N	O	S	0	0
			1608	495	820	146	144	3		

- Molecule 11 is a protein called Small ribosomal subunit protein uS11.

Mol	Chain	Residues	Atoms						AltConf	Trace
11	K	115	Total	C	H	N	O	S	0	0
			1719	528	864	170	156	1		

- Molecule 12 is a protein called Small ribosomal subunit protein uS12.

Mol	Chain	Residues	Atoms					AltConf	Trace	
12	L	122	Total	C	H	N	O	S	0	0
			2003	594	1045	197	165	2		

- Molecule 13 is a protein called Small ribosomal subunit protein uS13.

Mol	Chain	Residues	Atoms						AltConf	Trace
13	M	116	Total	C	H	N	O	S	0	0
			1922	572	987	191	169	3		

- Molecule 14 is a protein called Small ribosomal subunit protein uS14B.

Mol	Chain	Residues	Atoms						AltConf	Trace
14	N	60	Total	C	H	N	O	S	0	0
			981	302	504	97	73	5		

- Molecule 15 is a protein called Small ribosomal subunit protein uS15.

Mol	Chain	Residues	Atoms						AltConf	Trace
15	O	88	Total	C	H	N	O		0	0
			1481	449	761	147	124			

- Molecule 16 is a protein called Small ribosomal subunit protein bS16.

Mol	Chain	Residues	Atoms						AltConf	Trace
16	P	113	Total	C	H	N	O		0	0
			1827	570	936	162	159			

- Molecule 17 is a protein called Small ribosomal subunit protein uS17.

Mol	Chain	Residues	Atoms						AltConf	Trace
17	Q	94	Total	C	H	N	O	S	0	0
			1544	469	796	142	135	2		

- Molecule 18 is a protein called Small ribosomal subunit protein bS18B.

Mol	Chain	Residues	Atoms						AltConf	Trace
18	R	65	Total	C	H	N	O	S	0	0
			1054	318	541	102	90	3		

- Molecule 19 is a protein called Small ribosomal subunit protein uS19.

Mol	Chain	Residues	Atoms						AltConf	Trace
19	S	82	Total	C	H	N	O	S	0	0
			1339	425	677	124	112	1		

- Molecule 20 is a protein called Small ribosomal subunit protein bS20.

Mol	Chain	Residues	Atoms						AltConf	Trace
20	T	85	Total	C	H	N	O		0	0
			1373	402	713	139	119			

- Molecule 21 is a protein called Small ribosomal subunit protein uS2.

Mol	Chain	Residues	Atoms						AltConf	Trace
21	V	228	Total	C	H	N	O	S	0	0
			3632	1132	1839	322	330	9		

- Molecule 22 is a RNA chain called mRNA fragment.

Mol	Chain	Residues	Atoms						AltConf	Trace
22	X	6	Total	C	H	N	O	P	0	0
			180	54	63	13	45	5		

- Molecule 23 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
23	Y	2954	Total	C	H	N	O	P	0	0
			95352	28279	31908	11682	20529	2954		

- Molecule 24 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
24	U	118	Total	C	H	N	O	P	0	0
			3807	1126	1285	468	810	118		

- Molecule 25 is a protein called Large ribosomal subunit protein uL2.

Mol	Chain	Residues	Atoms						AltConf	Trace
25	Z	275	Total	C	H	N	O	S	0	0
			4276	1298	2166	438	370	4		

- Molecule 26 is a protein called Large ribosomal subunit protein uL3.

Mol	Chain	Residues	Atoms						AltConf	Trace
26	a	214	Total	C	H	N	O	S	0	0
			3218	982	1631	310	290	5		

- Molecule 27 is a protein called Large ribosomal subunit protein uL4.

Mol	Chain	Residues	Atoms						AltConf	Trace
27	b	209	Total	C	H	N	O	S	0	0
			3177	969	1608	295	303	2		

- Molecule 28 is a protein called Large ribosomal subunit protein uL5.



Mol	Chain	Residues	Atoms						AltConf	Trace
28	c	182	Total	C	H	N	O	S	0	0
			2922	907	1477	271	261	6		

- Molecule 29 is a protein called Large ribosomal subunit protein uL6.

Mol	Chain	Residues	Atoms						AltConf	Trace
29	d	176	Total	C	H	N	O	S	0	0
			2748	845	1400	249	253	1		

- Molecule 30 is a protein called Large ribosomal subunit protein uL11.

Mol	Chain	Residues	Atoms						AltConf	Trace
30	e	133	Total	C	H	N	O	S	0	0
			2012	625	1022	175	187	3		

- Molecule 31 is a protein called Large ribosomal subunit protein uL13.

Mol	Chain	Residues	Atoms						AltConf	Trace
31	f	146	Total	C	H	N	O	S	0	0
			2297	722	1167	207	200	1		

- Molecule 32 is a protein called Large ribosomal subunit protein uL14.

Mol	Chain	Residues	Atoms						AltConf	Trace
32	g	122	Total	C	H	N	O	S	0	0
			1938	586	1000	179	170	3		

- Molecule 33 is a protein called Large ribosomal subunit protein uL15.

Mol	Chain	Residues	Atoms						AltConf	Trace
33	h	145	Total	C	H	N	O	S	0	0
			2230	676	1152	205	194	3		

- Molecule 34 is a protein called Large ribosomal subunit protein uL16.

Mol	Chain	Residues	Atoms						AltConf	Trace
34	i	136	Total	C	H	N	O	S	0	0
			2220	690	1128	213	187	2		

- Molecule 35 is a protein called Large ribosomal subunit protein bL17.

Mol	Chain	Residues	Atoms						AltConf	Trace
35	j	118	Total	C	H	N	O	S	0	0
			1900	583	972	180	163	2		

- Molecule 36 is a protein called Large ribosomal subunit protein uL18.

Mol	Chain	Residues	Atoms						AltConf	Trace
36	k	126	Total	C	H	N	O		0	0
			1948	586	992	199	171			

- Molecule 37 is a protein called Large ribosomal subunit protein bL19.

Mol	Chain	Residues	Atoms						AltConf	Trace
37	l	113	Total	C	H	N	O	S	0	0
			1845	570	938	171	165	1		

- Molecule 38 is a protein called Large ribosomal subunit protein bL20.

Mol	Chain	Residues	Atoms						AltConf	Trace
38	m	124	Total	C	H	N	O		0	0
			2027	613	1039	203	172			

- Molecule 39 is a protein called Large ribosomal subunit protein bL21.

Mol	Chain	Residues	Atoms						AltConf	Trace
39	n	100	Total	C	H	N	O		0	0
			1557	478	803	137	139			

- Molecule 40 is a protein called Large ribosomal subunit protein uL22.

Mol	Chain	Residues	Atoms						AltConf	Trace
40	o	114	Total	C	H	N	O		0	0
			1783	543	910	171	159			

- Molecule 41 is a protein called Large ribosomal subunit protein uL23.

Mol	Chain	Residues	Atoms						AltConf	Trace
41	p	97	Total	C	H	N	O		0	0
			1559	479	803	138	139			

- Molecule 42 is a protein called Large ribosomal subunit protein uL24.

Mol	Chain	Residues	Atoms						AltConf	Trace
42	q	97	Total	C	H	N	O	S	0	0
			1515	456	783	137	137	2		

- Molecule 43 is a protein called Large ribosomal subunit protein bL25.

Mol	Chain	Residues	Atoms						AltConf	Trace
43	r	192	Total	C	H	N	O		0	0
			2872	881	1444	255	292			

- Molecule 44 is a protein called Large ribosomal subunit protein bL27.

Mol	Chain	Residues	Atoms						AltConf	Trace
44	s	79	Total	C	H	N	O		0	0
			1188	361	602	123	102			

- Molecule 45 is a protein called Large ribosomal subunit protein bL28.

Mol	Chain	Residues	Atoms						AltConf	Trace
45	t	63	Total	C	H	N	O	S	0	0
			955	283	485	103	80	4		

- Molecule 46 is a protein called Large ribosomal subunit protein uL29.

Mol	Chain	Residues	Atoms						AltConf	Trace
46	u	64	Total	C	H	N	O	S	0	0
			1073	324	542	103	103	1		

- Molecule 47 is a protein called 50S ribosomal protein bL37.

Mol	Chain	Residues	Atoms						AltConf	Trace
47	5	23	Total	C	H	N	O		0	0
			395	111	206	50	28			

- Molecule 48 is a protein called Large ribosomal subunit protein uL10.

Mol	Chain	Residues	Atoms						AltConf	Trace
48	v	126	Total	C	H	N	O	S	0	0
			1877	580	959	156	180	2		

- Molecule 49 is a protein called Large ribosomal subunit protein uL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	w	59	Total	C	H	N	O	0	0
			975	292	501	95	87		

- Molecule 50 is a protein called Large ribosomal subunit protein bL32.

Mol	Chain	Residues	Atoms						AltConf	Trace
50	x	54	Total	C	H	N	O	S	0	0
			887	260	464	93	69	1		

- Molecule 51 is a protein called Large ribosomal subunit protein bL33A.

Mol	Chain	Residues	Atoms						AltConf	Trace
51	y	49	Total	C	H	N	O	S	0	0
			817	248	412	82	71	4		

- Molecule 52 is a protein called Large ribosomal subunit protein bL34.

Mol	Chain	Residues	Atoms					AltConf	Trace	
52	z	46	Total	C	H	N	O	S	0	0
			789	225	412	97	54	1		

- Molecule 53 is a protein called Large ribosomal subunit protein bL35.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	1	63	Total	C	H	N	O	0	0
			1043	302	541	115	85		

- Molecule 54 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace	
54	2	37	Total	C	H	N	O	S	0	0
			623	181	324	66	47	5		

- Molecule 55 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
55	A	212	Total	Mg	0
			212	212	
55	B	1	Total	Mg	0
			1	1	
55	F	1	Total	Mg	0
			1	1	

*Continued on next page...*

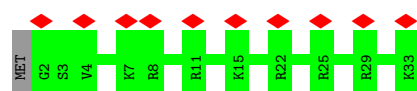
*Continued from previous page...*

Mol	Chain	Residues	Atoms		AltConf
55	N	1	Total 1	Mg 1	0
55	P	1	Total 1	Mg 1	0
55	R	1	Total 1	Mg 1	0
55	Y	381	Total 381	Mg 381	0
55	U	9	Total 9	Mg 9	0
55	Z	9	Total 9	Mg 9	0
55	b	1	Total 1	Mg 1	0
55	c	1	Total 1	Mg 1	0
55	i	1	Total 1	Mg 1	0
55	o	1	Total 1	Mg 1	0
55	p	1	Total 1	Mg 1	0
55	s	1	Total 1	Mg 1	0
55	t	1	Total 1	Mg 1	0
55	1	1	Total 1	Mg 1	0

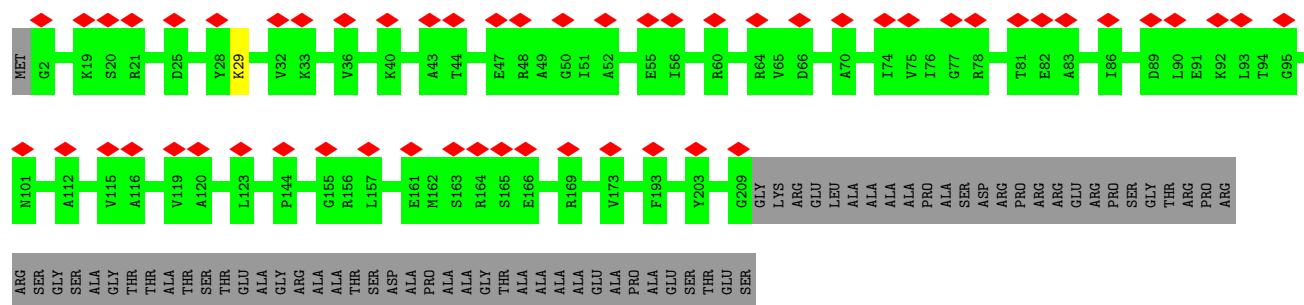
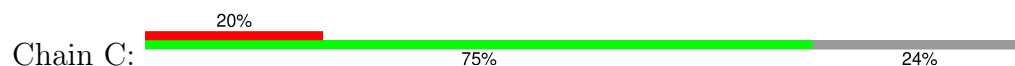
- Molecule 56 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
56	N	1	Total 1	Zn 1	0
56	R	1	Total 1	Zn 1	0
56	t	1	Total 1	Zn 1	0
56	y	1	Total 1	Zn 1	0
56	2	1	Total 1	Zn 1	0

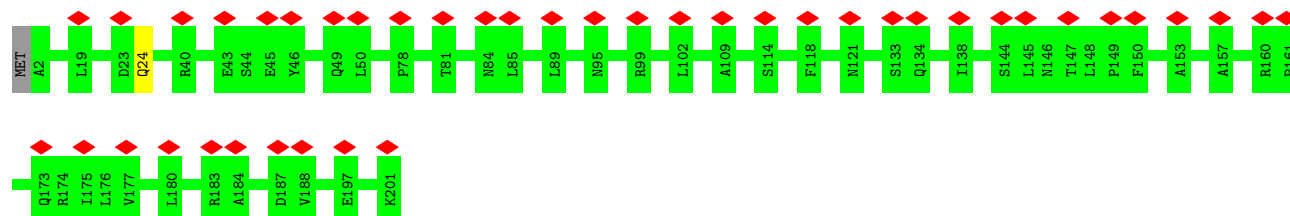




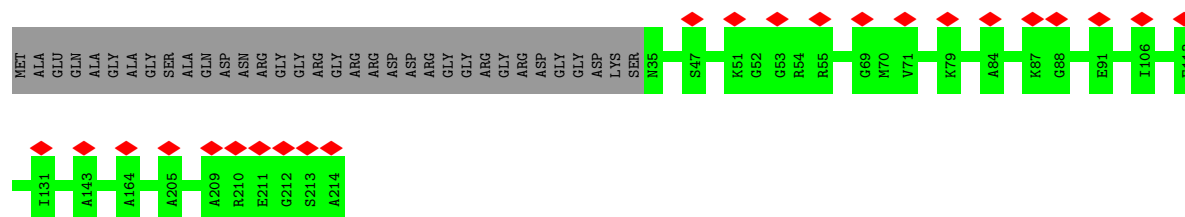
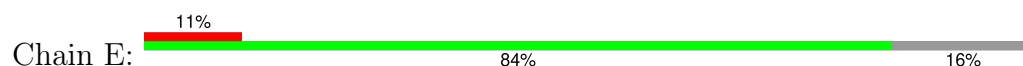
- Molecule 3: Small ribosomal subunit protein uS3



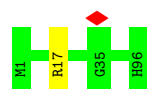
- Molecule 4: Small ribosomal subunit protein uS4



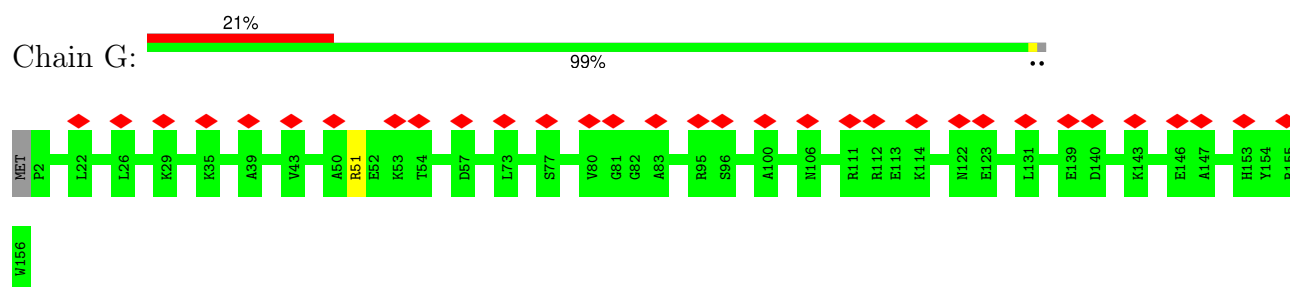
- Molecule 5: Small ribosomal subunit protein uS5



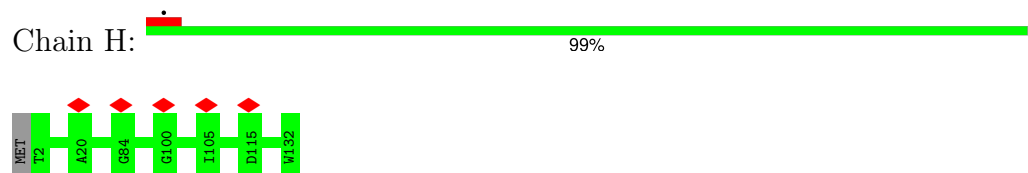
- Molecule 6: Small ribosomal subunit protein bS6



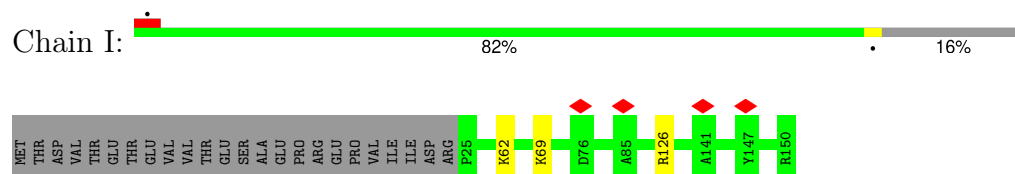
- Molecule 7: Small ribosomal subunit protein uS7



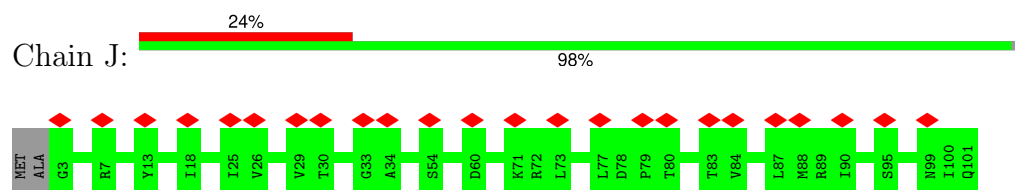
- Molecule 8: Small ribosomal subunit protein uS8



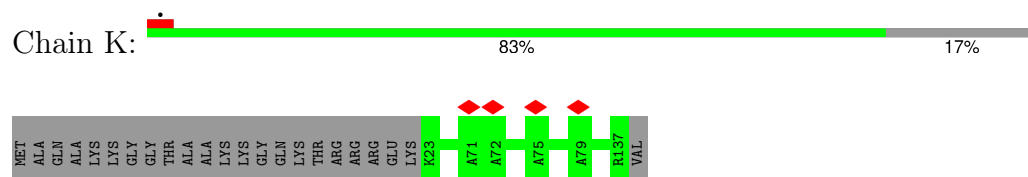
- Molecule 9: Small ribosomal subunit protein uS9



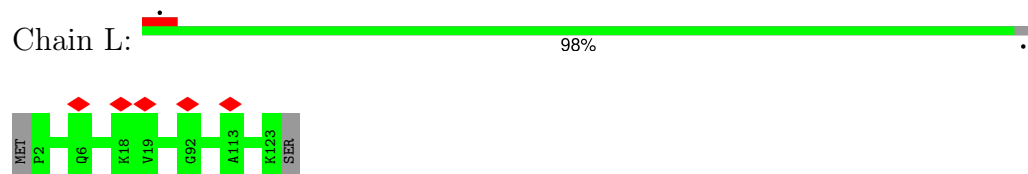
- Molecule 10: Small ribosomal subunit protein uS10



- Molecule 11: Small ribosomal subunit protein uS11



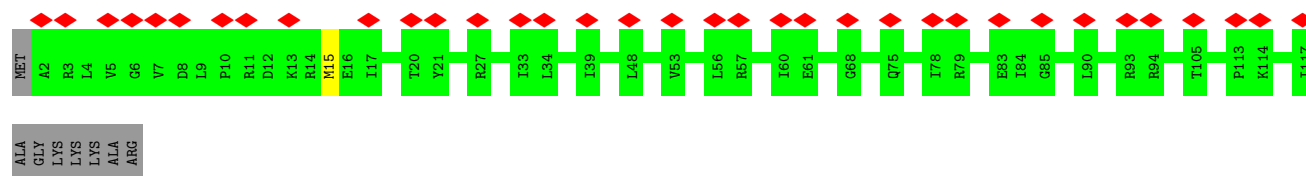
- Molecule 12: Small ribosomal subunit protein uS12



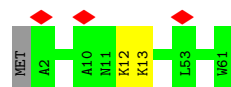
- Molecule 13: Small ribosomal subunit protein uS13



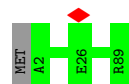




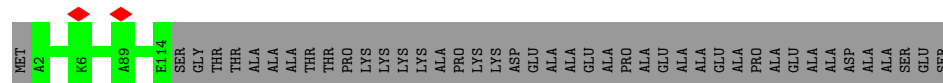
- Molecule 14: Small ribosomal subunit protein uS14B



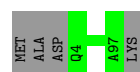
- Molecule 15: Small ribosomal subunit protein uS15



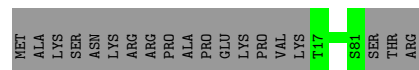
- Molecule 16: Small ribosomal subunit protein bS16



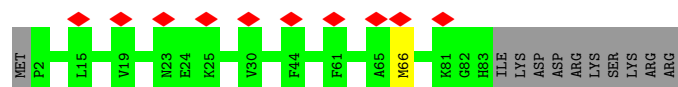
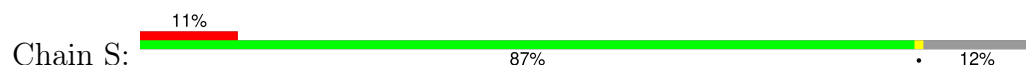
- Molecule 17: Small ribosomal subunit protein uS17



- Molecule 18: Small ribosomal subunit protein bS18B



- Molecule 19: Small ribosomal subunit protein uS19



- 

- |      |      |      |     |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |     |
|------|------|------|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|
| PRO  | ALA  | GLU  | GLY | ALA  | GLU  | PRO  | GLU  | ALA  | GLU  | TRP  | GLU  | GLN  | LEU  | LEU  | LEU  | ALA  | GLY  | THR  | ALA  | GLY  | ALA  | ALA  | ASP  | SER  | SER  | SER  | THR  | ASP  | ALA  | SER  |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |     |
| Q155 | K156 | V157 |     | V161 | W162 | V163 | V164 | D165 | T166 | N167 | K168 | E169 | H170 | V173 | G174 | R177 | K178 | T181 | P182 | V183 | L187 | D188 | T189 | N190 | D192 | P193 | D194 | V195 | V196 | D197 | Y198 | P199 | G202 | D205 | A206 | T207 | R208 | E223 | G224 | L225 | Q226 | A227 | R228 | ALA  | GLY  | GLN  | GLY  | SER  | GLY  | GLU  | LYS |
| E78  | R86  | M89  | Q94 | R95  | W96  | L97  | G98  | G99  | M100 | L101 | T102 | N103 | T106 | L111 | Q112 | R113 | L114 | K115 | E116 | L117 | E118 | A119 | M120 | E121 | Q122 | T123 | G124 | G125 | F126 | E127 | G128 | R129 | K132 | E133 | I134 | L135 | M136 | L137 | T138 | R139 | N142 | K143 | R146 | S147 | L148 | G149 | M150 | I151 | R152 | D153 |     |
| M1   | V4   | T5   | M6  | K7   | Q8   | L9   | L10  | D11  | S12  | G13  | A14  | H15  | F16  | G17  | H18  | Q19  | T20  | R21  | R22  | W23  | M27  | K28  | I31  | F32  | T33  | D34  | R35  | N36  | I40  | I41  | D42  | L43  | Q44  | Q45  | Y49  | I50  | A53  | V57  | K58  | V61  | A62  | H63  | G64  | F69  | T72  | Q75  |      |      |      |      |     |

- There are no outlier residues recorded for this chain.

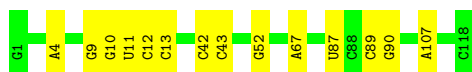
- |   |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |      |
|---|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|
| U | A2   | U7   | G20  | G21  | G23  | U31  | G32  | A60  | G61  | A68  | A71  | G72  | C73  | G78  | G82  | A89  | C90  | G94  | C95  | G96  | U97  | U98  | G99  | A115 | A116 | U117 | C125 | C126 | A148 | C172 | A180 | A189 | A195 | A212 | A215 | A220 | A221 |      |      |      |      |      |      |      |
|   | A227 | A228 | U229 | G230 | U231 | G248 | A272 | A273 | C274 | C275 | U279 | A282 | U285 | G286 | C291 | A296 | G300 | U301 | G302 | G303 | G311 | G312 | G313 | G314 | U315 | U316 | G317 | U318 | G319 | A322 | C323 | C324 | U330 | U331 | G335 | C336 | U337 | C338 | G342 | U343 | G344 | G352 | A355 | G356 |

A2826	C2646	G	U195	G1802	U1627	A1415	G1207	C1045	A667
U2833	U2647	U	G2196	A1803	A1628	A1416	U1208	C1046	A678
U2837	C2648	G	A2213	C1823	G1629	G1434	G1209	A1047	G679
C2853	A2649	A	C2214	C1824	A1631	C1435	C1210	G1048	G684
U2871	G2652	C	U2215	C1825	C1634	A1436	A1213	G1049	G685
A2814	A2653	A	G2216	A1834	A1635	A1437	U1219	G1063	A696
C2915	U2654	C	U2217	C1835	A1636	C1448	C1220	U1075	G706
A2926	U2655	C	A2221	G1849	G1637	C1449	A1077	A1076	G707
C2936	C2665	U	G2236	U2086	A1640	C1465	G1078	G1078	G708
G2937	G2671	G	A2237	U1864	U1641	G1473	C1222	A897	U709
C2938	A2672	A	C2247	A1865	C1649	A1474	U1226	U1084	G710
G2938	A2693	G	A2254	C1866	C1649	A1474	G1230	G1085	U714
C2938	G2694	U	A2255	U1870	G1674	A1480	U1231	U905	A721
U2953	C2698	C	A2257	G1871	U1675	A1493	A1233	A919	A731
U2953	C2699	A	A2258	A1872	G1676	U1494	U1234	G920	A738
A2957	A2700	C	C2267	G1892	A1679	A1499	G1240	G927	U739
A2972	U2701	C	G2096	C1912	U1681	A1510	G1251	A934	A740
A2973	A2702	G	A2106	A1916	G1703	U1511	G1252	U942	A747
A2974	G2705	C	G2107	G1917	U1784	G1522	C1253	G960	C749
G2975	G2713	G	U2111	G1933	C1705	C1531	C1260	U961	A756
C2976	G2714	U	U2112	U1946	A1710	G1532	A1261	U962	G757
A2981	U2715	G	G2130	U1947	U1713	U1533	A1262	G974	A758
A2982	C2722	U	G2131	G1950	A1714	C1534	G1270	U975	G759
G2985	G2726	G	A2137	C1973	U1717	U1540	A1274	C980	U760
G2986	A2727	U	C2138	A1974	U1728	G1541	U1292	U981	U764
A3002	U2728	G	U2139	A1975	A1729	A1542	G1293	A982	G765
C3008	G2729	U	A2140	U1981	A1731	A	C1298	U995	G766
U3009	U2737	C	U2141	A1990	A1737	U	G1301	G996	U767
C3013	A2742	G	A2143	A1998	G1746	C	U1341	G997	G768
A3014	U2743	C	C2148	C1998	U1754	C	U1344	G998	U769
C3015	C2744	U	U2149	A2001	A1755	A	A1344	C1002	G784
A3021	G2753	U	G2150	A2008	G1756	C	G1350	A1003	U801
G3022	A2790	U	C2153	A2008	U1757	G	G1353	C1004	G828
A3042	G2791	G	G2154	C2017	G1758	U	A1191	A1005	U829
G3070	A2796	A	A2162	G2018	U1767	A	G1371	U1010	A830
C3088	G2806	U	U2163	C2025	U1786	C	U1379	A1011	A831
A3093	U2809	C	C2166	A2026	A1787	C	G1384	U1012	U839
U2820	G2809	A	U2167	U2033	G1788	C	G1385	G1013	G840
C2821	G2809	C	U2179	G2034	A1789	C	U1387	G1014	G841
A2822	U2822	A	U2187	U2835	A1789	C	A1387	A1025	C845
C3105	G2823	U	C2191	A2036	U1798	A	U1389	A1032	U862
C3106	A2630	U	A2194	A2037	C1801	A	U1389	A1033	G863
				A2038				U1034	



- Molecule 24: 5S rRNA

Chain U: 88% 12%



- Molecule 25: Large ribosomal subunit protein uL2

Chain Z: 99%



- Molecule 26: Large ribosomal subunit protein uL3

Chain a: 98%



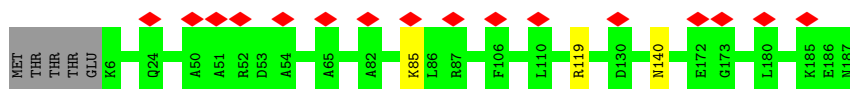
- Molecule 27: Large ribosomal subunit protein uL4

Chain b: 96%



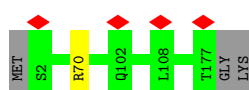
- Molecule 28: Large ribosomal subunit protein uL5

Chain c: 9% 96%



- Molecule 29: Large ribosomal subunit protein uL6

Chain d: 98%

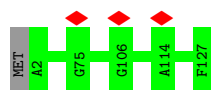


- Molecule 30: Large ribosomal subunit protein uL11



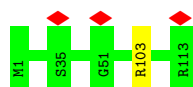
- Molecule 36: Large ribosomal subunit protein uL18

Chain k:  99%



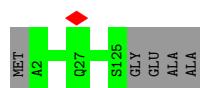
- Molecule 37: Large ribosomal subunit protein bL19

Chain l:  99%



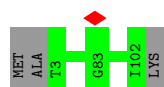
- Molecule 38: Large ribosomal subunit protein bL20

Chain m:  96%



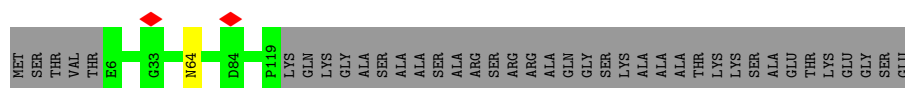
- Molecule 39: Large ribosomal subunit protein bL21

Chain n:  97%



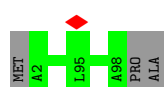
- Molecule 40: Large ribosomal subunit protein uL22

Chain o:  74% 25%

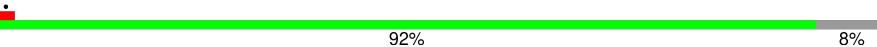


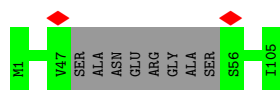
- Molecule 41: Large ribosomal subunit protein uL23

Chain p:  97%

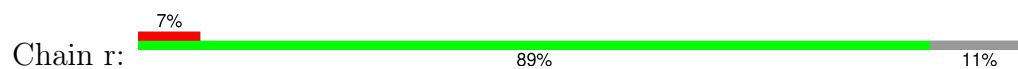


- Molecule 42: Large ribosomal subunit protein uL24

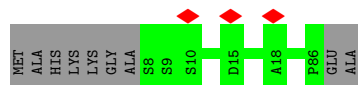
Chain q:  92% 8%



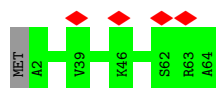
- Molecule 43: Large ribosomal subunit protein bL25



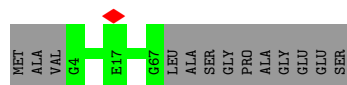
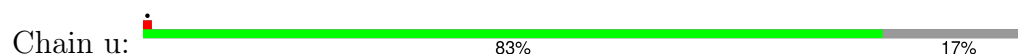
- Molecule 44: Large ribosomal subunit protein bL27



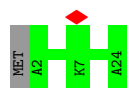
- Molecule 45: Large ribosomal subunit protein bL28



- Molecule 46: Large ribosomal subunit protein uL29



- Molecule 47: 50S ribosomal protein bL37



- Molecule 48: Large ribosomal subunit protein uL10



PRO  
ALA  
SER  
GLN  
VAL  
ALA  
ARG  
LEU  
ALA  
ALA  
ALA  
LEU  
GLN  
GLU  
LYS  
ALA  
GLY  
GLU  
GLU  
ALA  
ALA

- Molecule 49: Large ribosomal subunit protein uL30

Chain w:  97% .


MET  
A2  
G60  
LYS

- Molecule 50: Large ribosomal subunit protein bL32

Chain x:  95% 5%

MET  
A2  
D55  
LYS  
ARG

- Molecule 51: Large ribosomal subunit protein bL33A

Chain y:  89% 11%

MET  
ALA  
SER  
SER  
THR  
D6  
S54  
ARG

- Molecule 52: Large ribosomal subunit protein bL34

Chain z:  98% .

MET  
A2  
A47

- Molecule 53: Large ribosomal subunit protein bL35

Chain 1:  98% .

MET  
F2  
G64

- Molecule 54: 50S ribosomal protein L36

Chain 2:  5% 100%

H1  
G21  
G37



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	136418	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	72.704	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	130000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	17.694	Depositor
Minimum map value	-13.756	Depositor
Average map value	-0.010	Depositor
Map value standard deviation	0.546	Depositor
Recommended contour level	0.901	Depositor
Map size ( $\text{\AA}$ )	573.696, 573.696, 573.696	wwPDB
Map dimensions	864, 864, 864	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.664, 0.664, 0.664	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	A	0.21	0/36308	0.74	4/56653 (0.0%)
2	B	0.28	0/280	0.67	0/359
3	C	0.25	0/1684	0.56	0/2261
4	D	0.25	0/1672	0.55	0/2251
5	E	0.25	0/1312	0.54	0/1772
6	F	0.25	0/782	0.54	0/1059
7	G	0.23	0/1252	0.55	0/1690
8	H	0.25	0/1025	0.54	0/1385
9	I	0.25	0/1012	0.56	0/1362
10	J	0.26	0/802	0.54	0/1086
11	K	0.26	0/873	0.51	0/1180
12	L	0.24	0/969	0.60	0/1294
13	M	0.26	0/942	0.63	0/1260
14	N	0.26	0/488	0.55	0/650
15	O	0.25	0/729	0.58	0/977
16	P	0.26	0/908	0.54	0/1226
17	Q	0.25	0/759	0.56	0/1016
18	R	0.26	0/518	0.61	0/693
19	S	0.24	0/680	0.49	0/915
20	T	0.26	0/663	0.60	0/882
21	V	0.25	0/1822	0.52	0/2457
22	X	0.13	0/128	0.64	0/196
23	Y	0.25	0/71012	0.77	15/110722 (0.0%)
24	U	0.22	0/2821	0.74	0/4396
25	Z	0.26	0/2153	0.60	0/2895
26	a	0.25	0/1609	0.56	0/2165
27	b	0.25	0/1592	0.50	0/2153
28	c	0.26	0/1467	0.57	0/1973
29	d	0.25	0/1369	0.53	0/1848
30	e	0.24	0/986	0.44	0/1303
31	f	0.24	0/1157	0.48	0/1567
32	g	0.26	0/946	0.55	0/1268

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
33	h	0.25	0/1091	0.52	0/1457
34	i	0.25	0/1118	0.57	0/1506
35	j	0.26	0/945	0.54	0/1267
36	k	0.25	0/966	0.58	0/1298
37	l	0.26	0/921	0.56	0/1236
38	m	0.27	0/1000	0.55	0/1341
39	n	0.25	0/764	0.48	0/1030
40	o	0.25	0/887	0.56	0/1204
41	p	0.26	0/766	0.55	0/1030
42	q	0.24	0/738	0.50	0/987
43	r	0.24	0/1443	0.50	0/1970
44	s	0.26	0/595	0.57	0/798
45	t	0.25	0/478	0.59	0/641
46	u	0.29	0/534	0.64	0/713
47	5	0.24	0/191	0.62	0/247
48	v	0.25	0/925	0.47	0/1246
49	w	0.24	0/477	0.58	0/640
50	x	0.24	0/427	0.61	0/572
51	y	0.26	0/413	0.54	0/553
52	z	0.25	0/380	0.69	0/500
53	1	0.22	0/507	0.62	0/672
54	2	0.25	0/303	0.57	0/401
All	All	0.24	0/156589	0.71	19/234223 (0.0%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1382	C	OP2-P-O3'	-11.22	80.51	105.20
23	Y	72	G	OP1-P-O3'	-10.64	81.79	105.20
23	Y	72	G	OP2-P-O3'	-9.85	83.52	105.20
23	Y	1449	C	N3-C2-O2	-8.84	115.72	121.90
1	A	1382	C	OP1-P-O3'	-8.66	86.15	105.20

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	
2	B	30/33 (91%)	29 (97%)	1 (3%)	0	100	100
3	C	206/275 (75%)	197 (96%)	9 (4%)	0	100	100
4	D	198/201 (98%)	184 (93%)	14 (7%)	0	100	100
5	E	178/214 (83%)	172 (97%)	6 (3%)	0	100	100
6	F	94/96 (98%)	91 (97%)	3 (3%)	0	100	100
7	G	153/156 (98%)	152 (99%)	1 (1%)	0	100	100
8	H	129/132 (98%)	126 (98%)	3 (2%)	0	100	100
9	I	124/150 (83%)	112 (90%)	12 (10%)	0	100	100
10	J	97/101 (96%)	92 (95%)	5 (5%)	0	100	100
11	K	113/138 (82%)	106 (94%)	7 (6%)	0	100	100
12	L	120/124 (97%)	110 (92%)	10 (8%)	0	100	100
13	M	114/124 (92%)	110 (96%)	4 (4%)	0	100	100
14	N	58/61 (95%)	53 (91%)	5 (9%)	0	100	100
15	O	86/89 (97%)	85 (99%)	1 (1%)	0	100	100
16	P	111/156 (71%)	106 (96%)	5 (4%)	0	100	100
17	Q	92/98 (94%)	89 (97%)	3 (3%)	0	100	100
18	R	63/84 (75%)	59 (94%)	4 (6%)	0	100	100
19	S	80/93 (86%)	78 (98%)	2 (2%)	0	100	100
20	T	83/86 (96%)	83 (100%)	0	0	100	100
21	V	226/277 (82%)	218 (96%)	6 (3%)	2 (1%)	14	43
25	Z	273/278 (98%)	263 (96%)	9 (3%)	1 (0%)	30	60
26	a	212/217 (98%)	196 (92%)	16 (8%)	0	100	100
27	b	207/215 (96%)	200 (97%)	7 (3%)	0	100	100
28	c	180/187 (96%)	174 (97%)	6 (3%)	0	100	100
29	d	174/179 (97%)	169 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
30	e	96/142 (68%)	93 (97%)	3 (3%)	0	100	100
31	f	144/147 (98%)	139 (96%)	5 (4%)	0	100	100
32	g	120/122 (98%)	118 (98%)	2 (2%)	0	100	100
33	h	143/147 (97%)	129 (90%)	14 (10%)	0	100	100
34	i	134/138 (97%)	127 (95%)	7 (5%)	0	100	100
35	j	116/199 (58%)	107 (92%)	9 (8%)	0	100	100
36	k	124/127 (98%)	122 (98%)	2 (2%)	0	100	100
37	l	111/113 (98%)	104 (94%)	7 (6%)	0	100	100
38	m	122/129 (95%)	118 (97%)	4 (3%)	0	100	100
39	n	98/103 (95%)	95 (97%)	3 (3%)	0	100	100
40	o	112/153 (73%)	107 (96%)	5 (4%)	0	100	100
41	p	95/100 (95%)	92 (97%)	3 (3%)	0	100	100
42	q	93/105 (89%)	91 (98%)	2 (2%)	0	100	100
43	r	190/215 (88%)	184 (97%)	6 (3%)	0	100	100
44	s	77/88 (88%)	74 (96%)	3 (4%)	0	100	100
45	t	61/64 (95%)	58 (95%)	3 (5%)	0	100	100
46	u	62/77 (80%)	61 (98%)	1 (2%)	0	100	100
47	5	21/24 (88%)	20 (95%)	1 (5%)	0	100	100
48	v	124/175 (71%)	122 (98%)	1 (1%)	1 (1%)	16	45
49	w	57/61 (93%)	54 (95%)	3 (5%)	0	100	100
50	x	52/57 (91%)	52 (100%)	0	0	100	100
51	y	47/55 (86%)	45 (96%)	2 (4%)	0	100	100
52	z	44/47 (94%)	44 (100%)	0	0	100	100
53	1	61/64 (95%)	57 (93%)	4 (7%)	0	100	100
54	2	35/37 (95%)	33 (94%)	2 (6%)	0	100	100
All	All	5740/6453 (89%)	5500 (96%)	236 (4%)	4 (0%)	50	77

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
21	V	155	GLN
25	Z	221	VAL
48	v	85	GLU

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Mol	Chain	Res	Type
21	V	157	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	30/31 (97%)	30 (100%)	0	100	100
3	C	170/212 (80%)	169 (99%)	1 (1%)	84	90
4	D	175/176 (99%)	174 (99%)	1 (1%)	84	90
5	E	127/147 (86%)	127 (100%)	0	100	100
6	F	85/85 (100%)	84 (99%)	1 (1%)	67	82
7	G	131/132 (99%)	130 (99%)	1 (1%)	79	88
8	H	107/108 (99%)	107 (100%)	0	100	100
9	I	102/125 (82%)	99 (97%)	3 (3%)	37	62
10	J	89/90 (99%)	89 (100%)	0	100	100
11	K	89/105 (85%)	89 (100%)	0	100	100
12	L	103/105 (98%)	103 (100%)	0	100	100
13	M	99/104 (95%)	98 (99%)	1 (1%)	73	85
14	N	49/50 (98%)	47 (96%)	2 (4%)	26	54
15	O	76/77 (99%)	76 (100%)	0	100	100
16	P	92/118 (78%)	92 (100%)	0	100	100
17	Q	80/83 (96%)	80 (100%)	0	100	100
18	R	55/72 (76%)	55 (100%)	0	100	100
19	S	73/84 (87%)	72 (99%)	1 (1%)	62	79
20	T	69/70 (99%)	66 (96%)	3 (4%)	25	53
21	V	191/218 (88%)	189 (99%)	2 (1%)	73	85
25	Z	215/218 (99%)	215 (100%)	0	100	100
26	a	160/163 (98%)	159 (99%)	1 (1%)	84	90
27	b	169/173 (98%)	167 (99%)	2 (1%)	67	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
28	c	151/156 (97%)	148 (98%)	3 (2%)	50	71
29	d	148/150 (99%)	147 (99%)	1 (1%)	81	89
30	e	102/108 (94%)	102 (100%)	0	100	100
31	f	119/120 (99%)	119 (100%)	0	100	100
32	g	100/100 (100%)	100 (100%)	0	100	100
33	h	112/114 (98%)	111 (99%)	1 (1%)	75	86
34	i	114/116 (98%)	114 (100%)	0	100	100
35	j	97/158 (61%)	97 (100%)	0	100	100
36	k	93/94 (99%)	93 (100%)	0	100	100
37	l	100/100 (100%)	99 (99%)	1 (1%)	73	85
38	m	97/99 (98%)	97 (100%)	0	100	100
39	n	81/83 (98%)	81 (100%)	0	100	100
40	o	90/117 (77%)	89 (99%)	1 (1%)	70	83
41	p	83/85 (98%)	83 (100%)	0	100	100
42	q	81/86 (94%)	81 (100%)	0	100	100
43	r	155/168 (92%)	155 (100%)	0	100	100
44	s	58/63 (92%)	58 (100%)	0	100	100
45	t	50/51 (98%)	50 (100%)	0	100	100
46	u	58/66 (88%)	58 (100%)	0	100	100
47	5	18/19 (95%)	18 (100%)	0	100	100
48	v	89/120 (74%)	89 (100%)	0	100	100
49	w	52/54 (96%)	52 (100%)	0	100	100
50	x	43/46 (94%)	43 (100%)	0	100	100
51	y	47/52 (90%)	47 (100%)	0	100	100
52	z	35/36 (97%)	35 (100%)	0	100	100
53	1	53/54 (98%)	53 (100%)	0	100	100
54	2	35/35 (100%)	35 (100%)	0	100	100
All	All	4797/5196 (92%)	4771 (100%)	26 (0%)	85	91

5 of 26 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
21	V	86	ARG

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Mol	Chain	Res	Type
27	b	83	GLN
37	l	103	ARG
26	a	159	ARG
27	b	171	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
4	D	146	ASN
9	I	66	GLN
9	I	146	GLN
25	Z	129	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1509/1528 (98%)	292 (19%)	16 (1%)
22	X	5/6 (83%)	0	0
23	Y	2921/3120 (93%)	530 (18%)	34 (1%)
24	U	117/118 (99%)	13 (11%)	1 (0%)
All	All	4552/4772 (95%)	835 (18%)	51 (1%)

5 of 835 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	8	U
1	A	9	U
1	A	11	G
1	A	12	A
1	A	13	G

5 of 51 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
23	Y	981	U
23	Y	1436	C
23	Y	3008	C
23	Y	1004	C
23	Y	1084	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 629 ligands modelled in this entry, 629 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
23	Y	31
30	e	20
1	A	1

The worst 5 of 52 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	Y	2406:U	O3'	2407:C	P	20.82
1	Y	1008:G	O3'	1009:U	P	11.65
1	e	22:PRO	C	23:ALA	N	10.56
1	Y	2096:G	O3'	2097:G	P	8.31
1	e	15:ILE	C	16:GLN	N	8.16

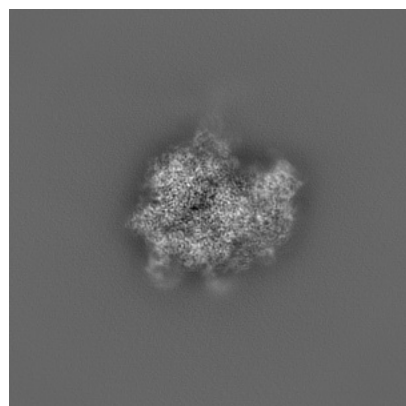
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44091. These allow visual inspection of the internal detail of the map and identification of artifacts.

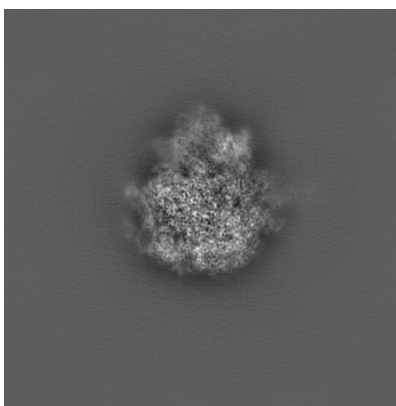
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

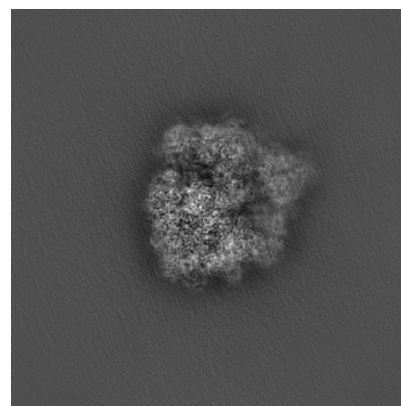
#### 6.1.1 Primary map



X

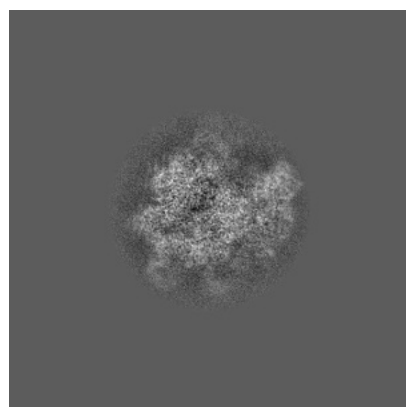


Y

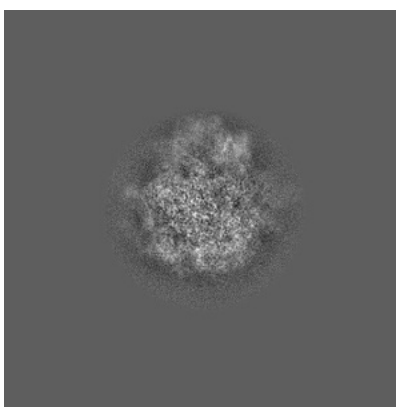


Z

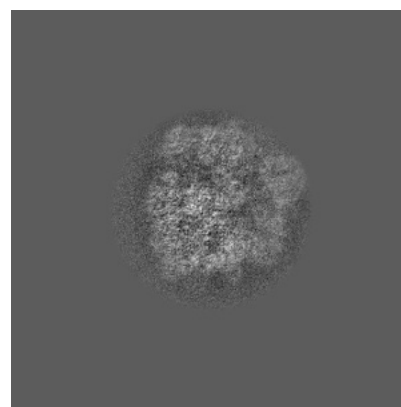
#### 6.1.2 Raw map



X



Y

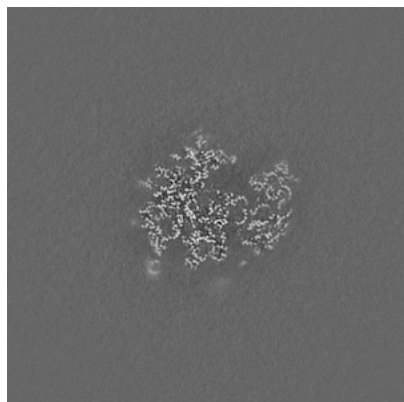


Z

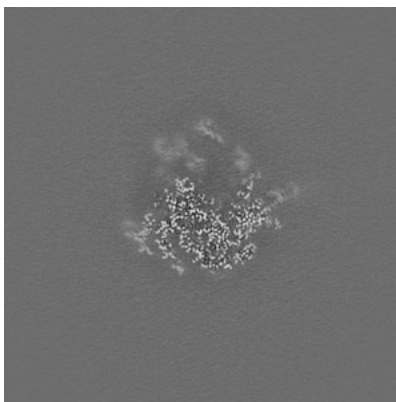
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

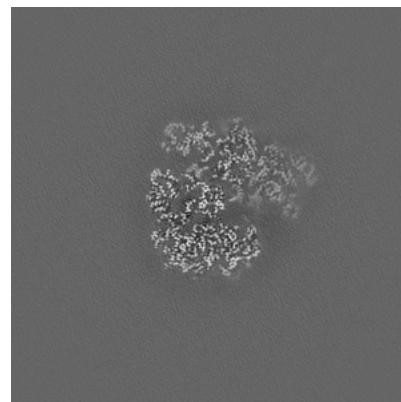
### 6.2.1 Primary map



X Index: 432

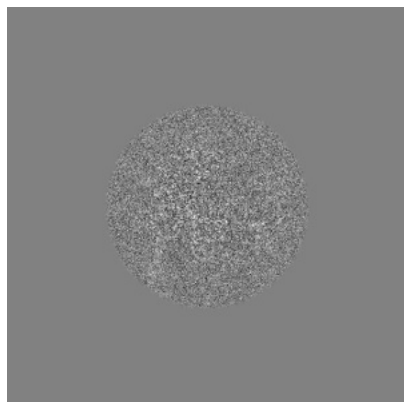


Y Index: 432

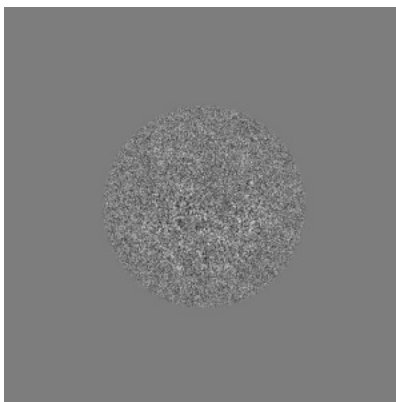


Z Index: 432

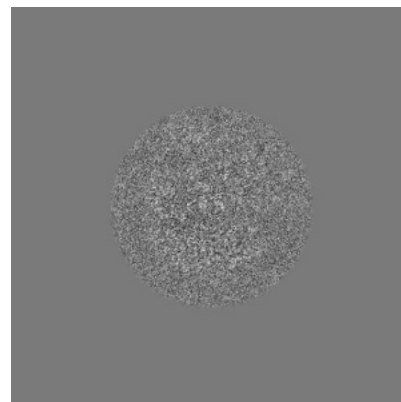
### 6.2.2 Raw map



X Index: 432



Y Index: 432

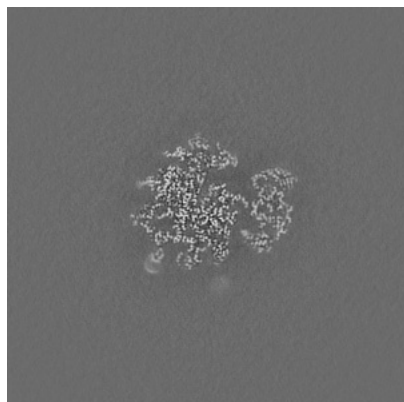


Z Index: 432

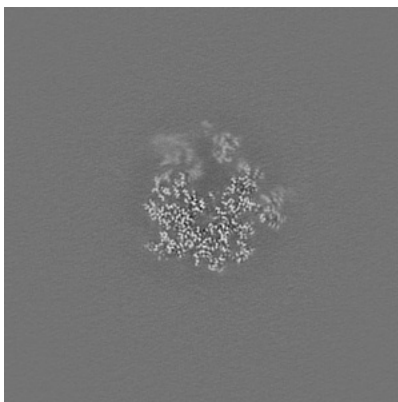
The images above show central slices of the map in three orthogonal directions.

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

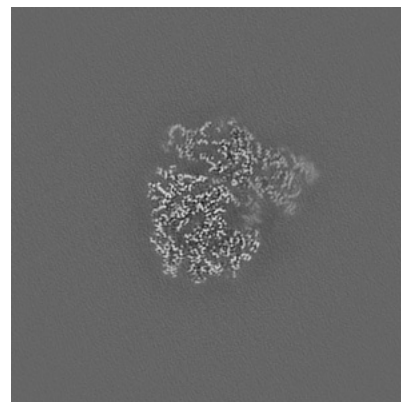
### 6.3.1 Primary map



X Index: 423

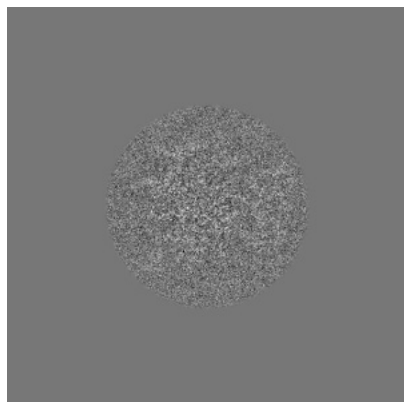


Y Index: 411

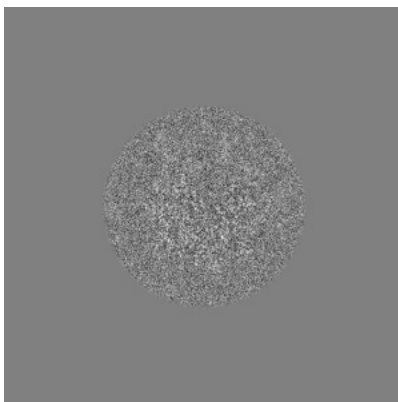


Z Index: 423

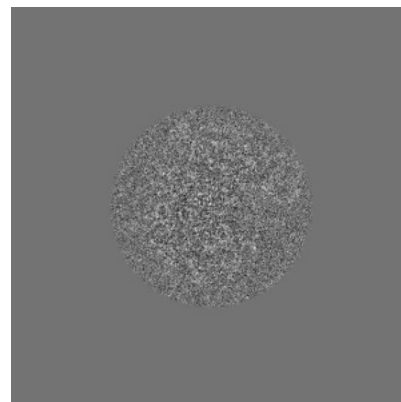
### 6.3.2 Raw map



X Index: 423



Y Index: 411

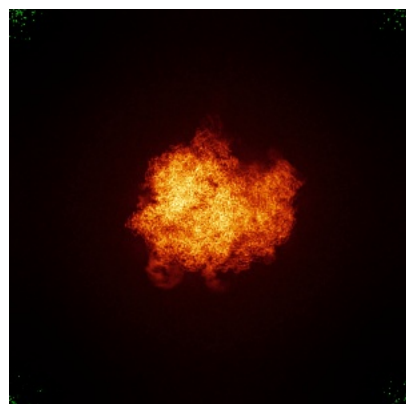


Z Index: 417

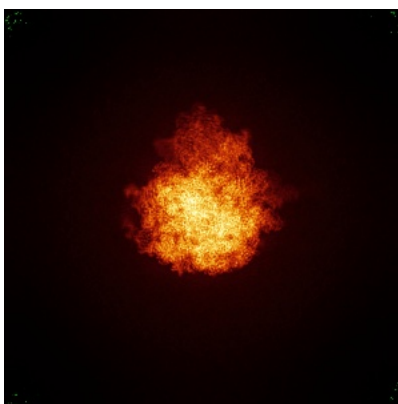
The images above show the largest variance slices of the map in three orthogonal directions.

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

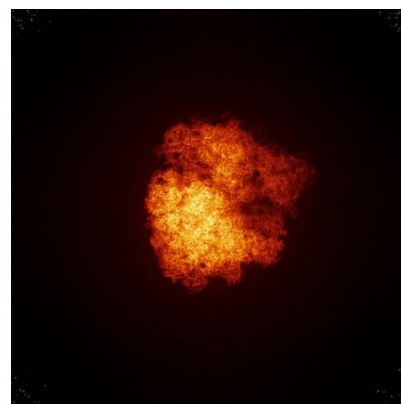
### 6.4.1 Primary map



X

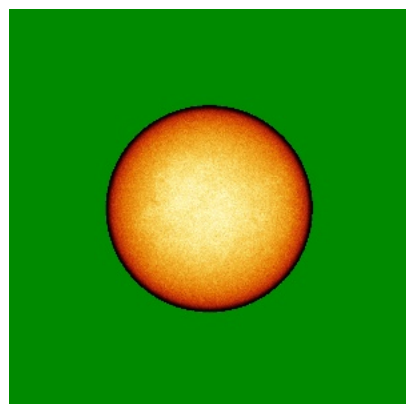


Y

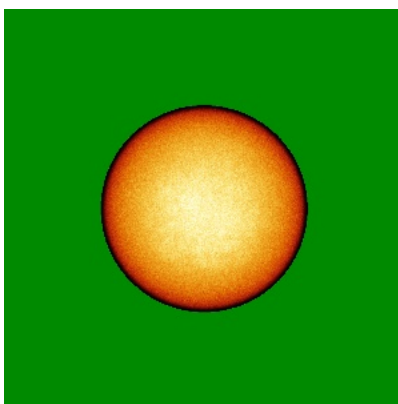


Z

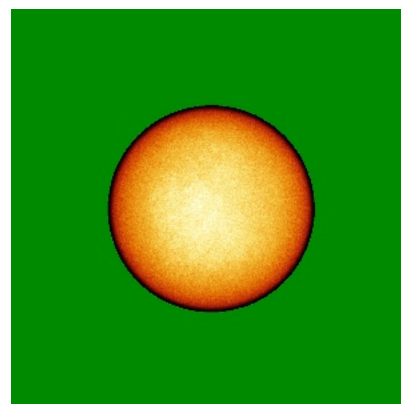
### 6.4.2 Raw map



X



Y

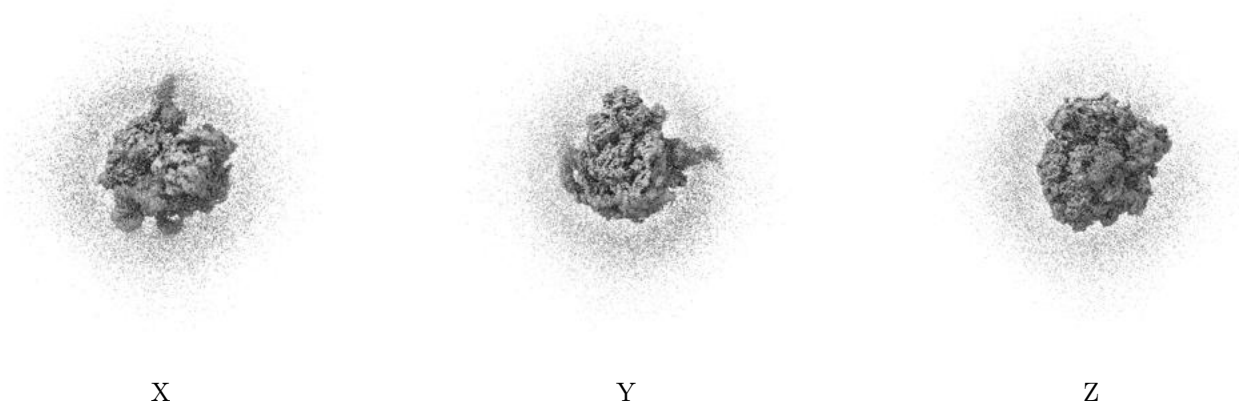


Z

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.

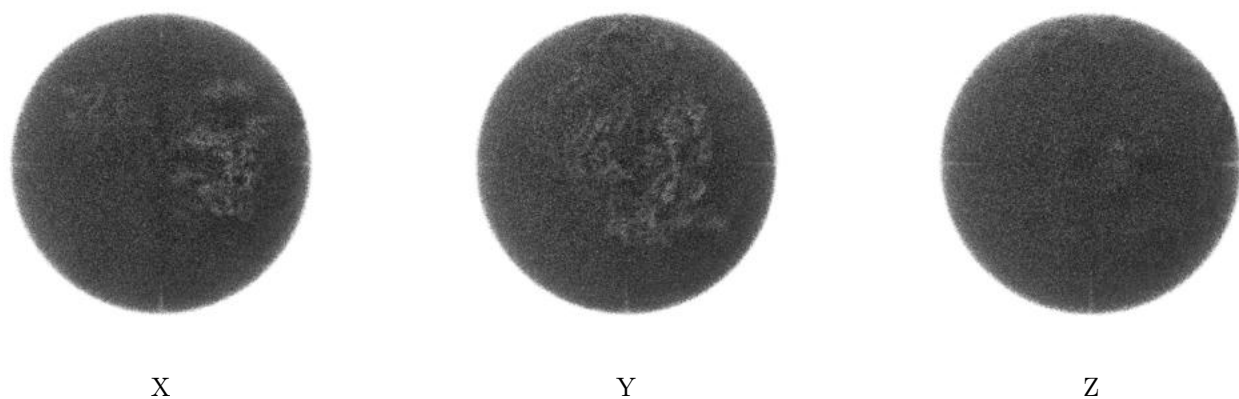
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.901. 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.

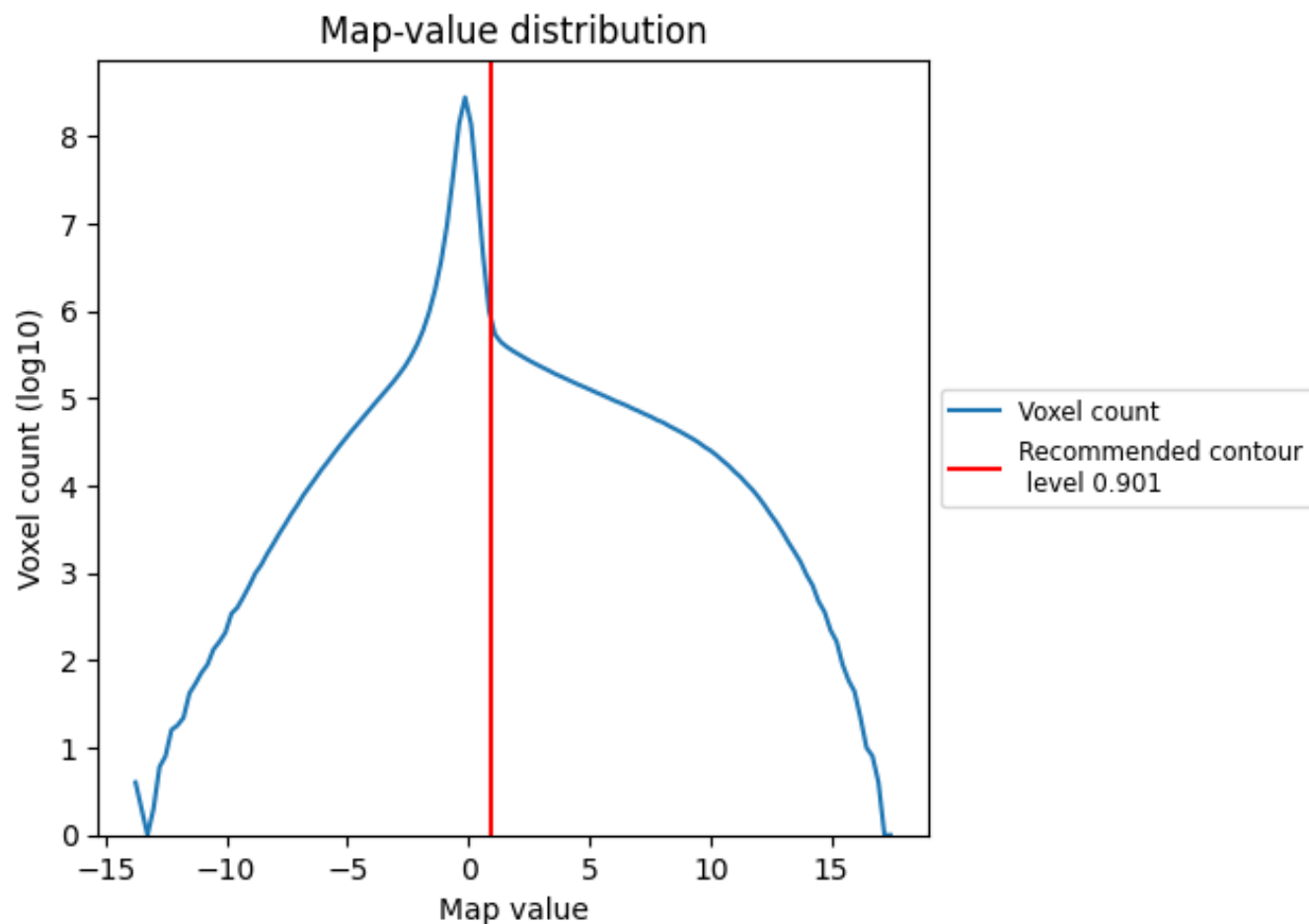
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

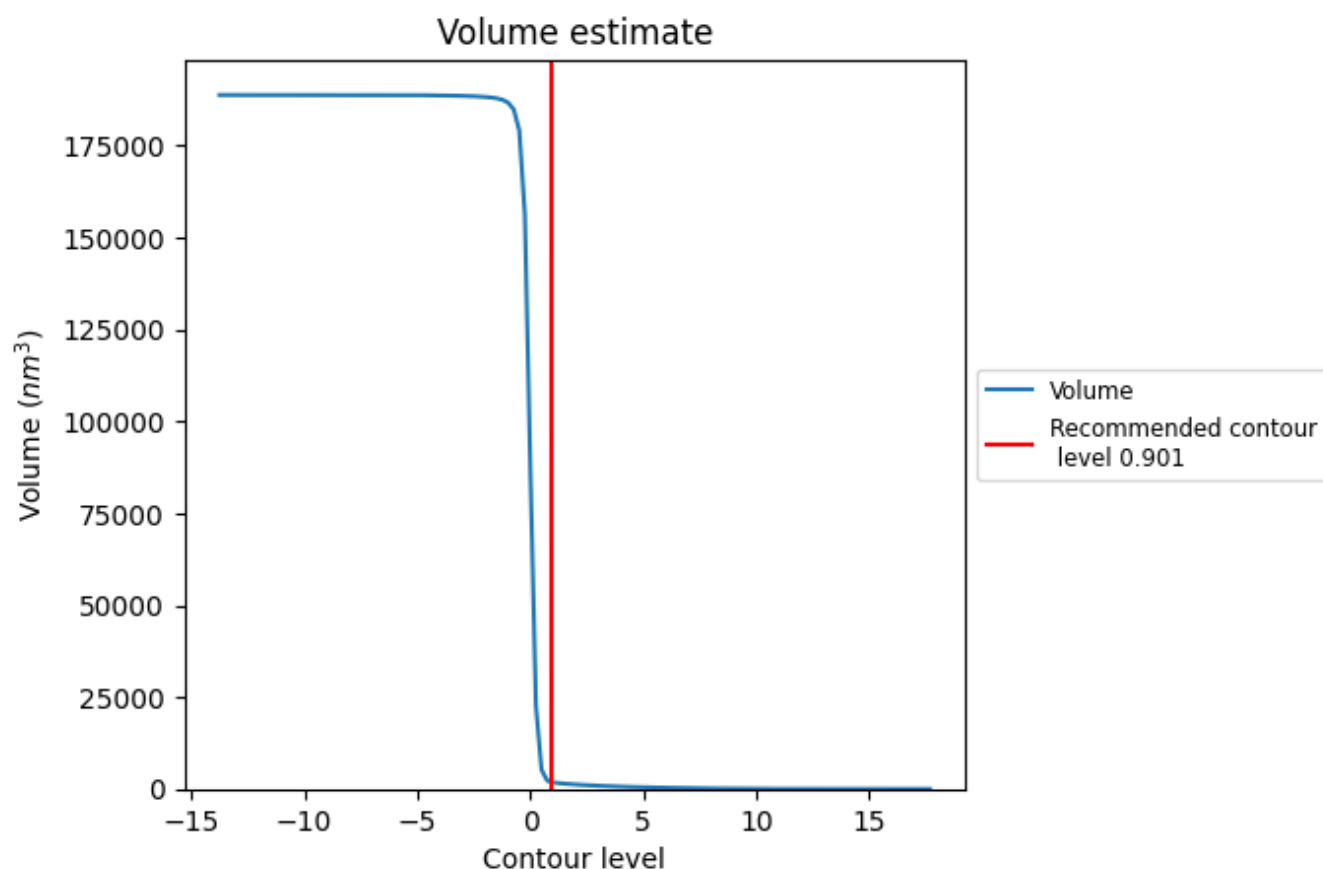
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

## 7.2 Volume estimate [i](#)

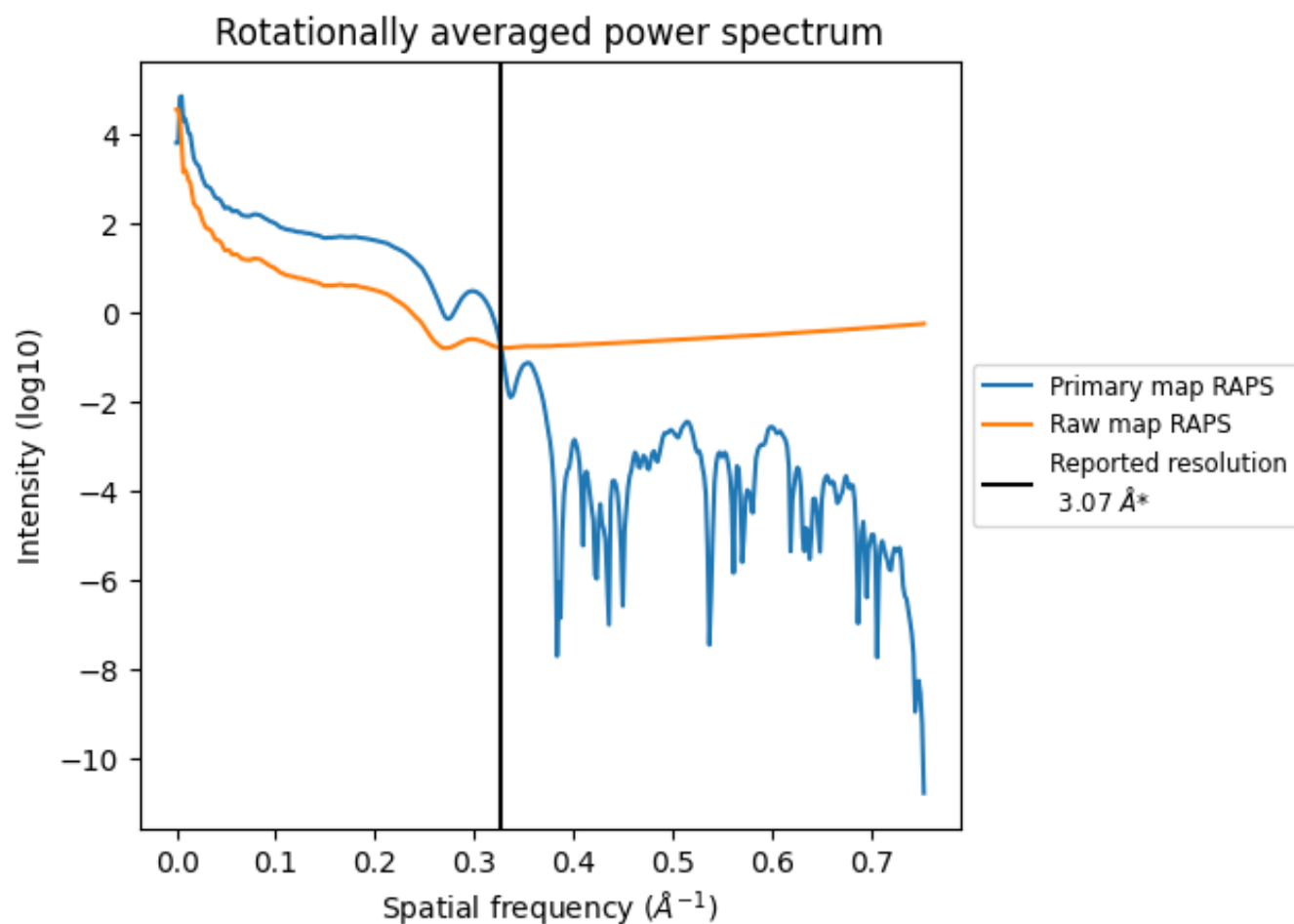


The volume at the recommended contour level is 1958  $\text{nm}^3$ ; this corresponds to an approximate mass of 1768 kDa.

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 ⓘ

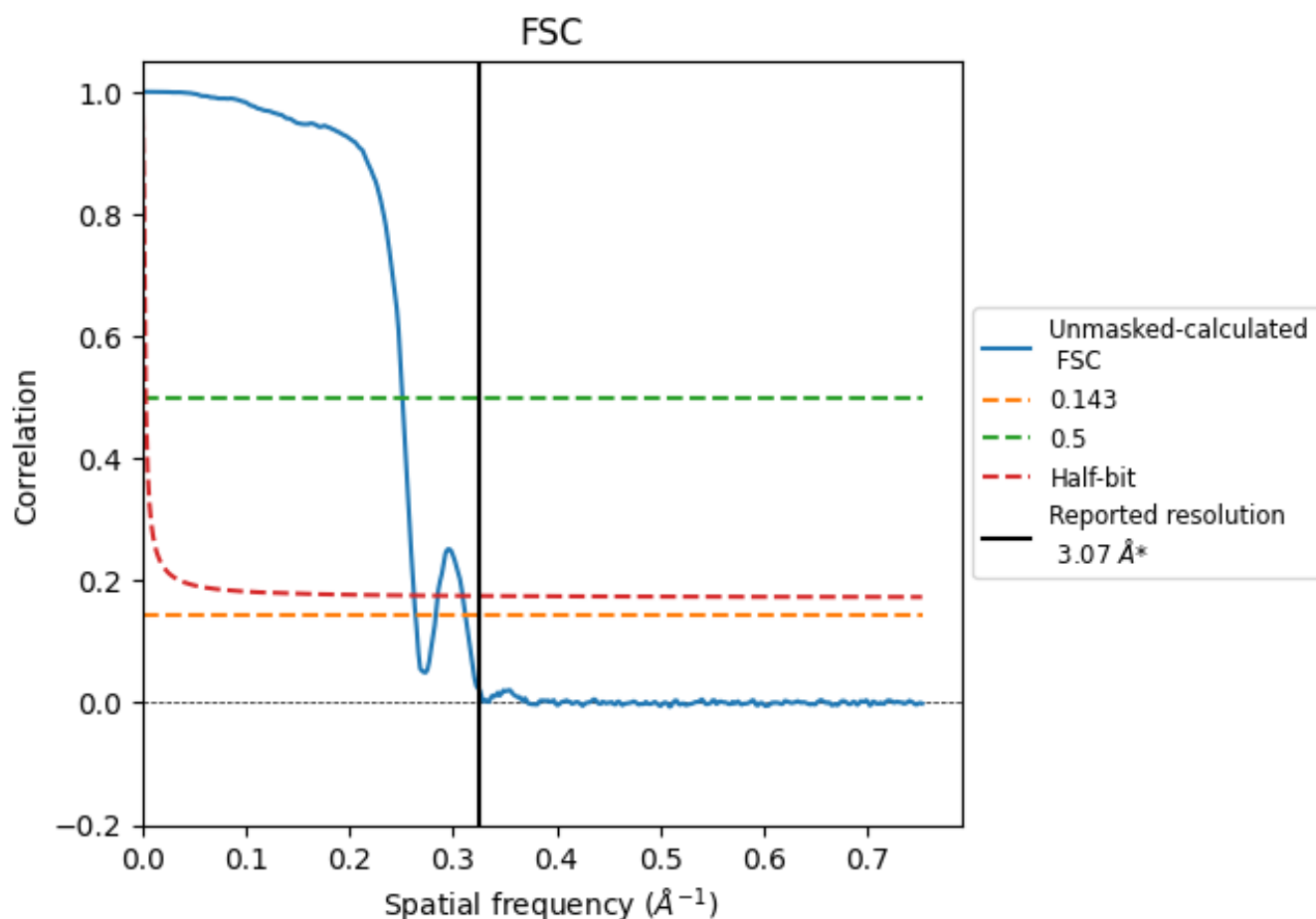


\*Reported resolution corresponds to spatial frequency of 0.326 Å<sup>-1</sup>

## 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 0.326 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

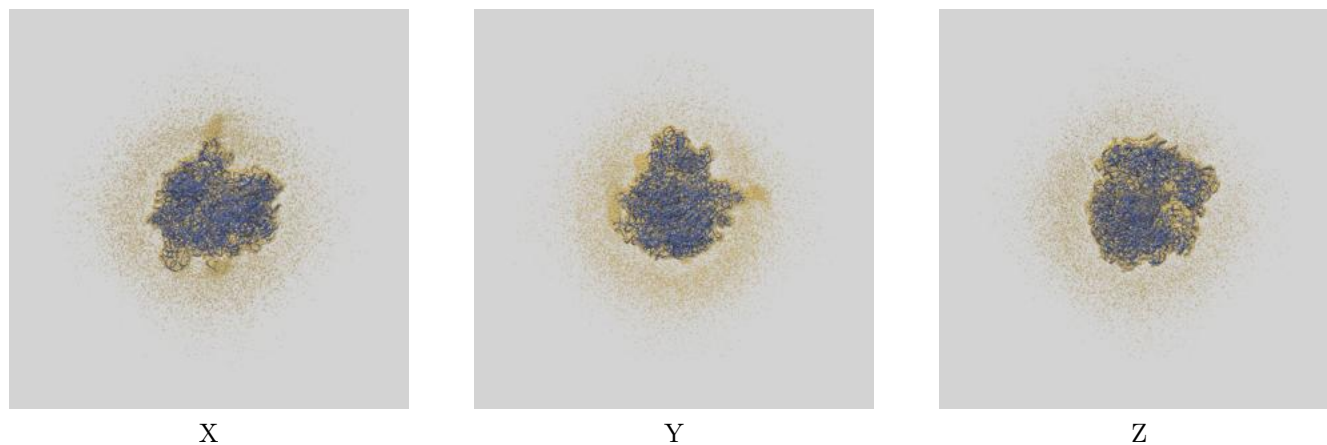
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.07	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.79	3.98	3.81

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.79 differs from the reported value 3.07 by more than 10 %

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

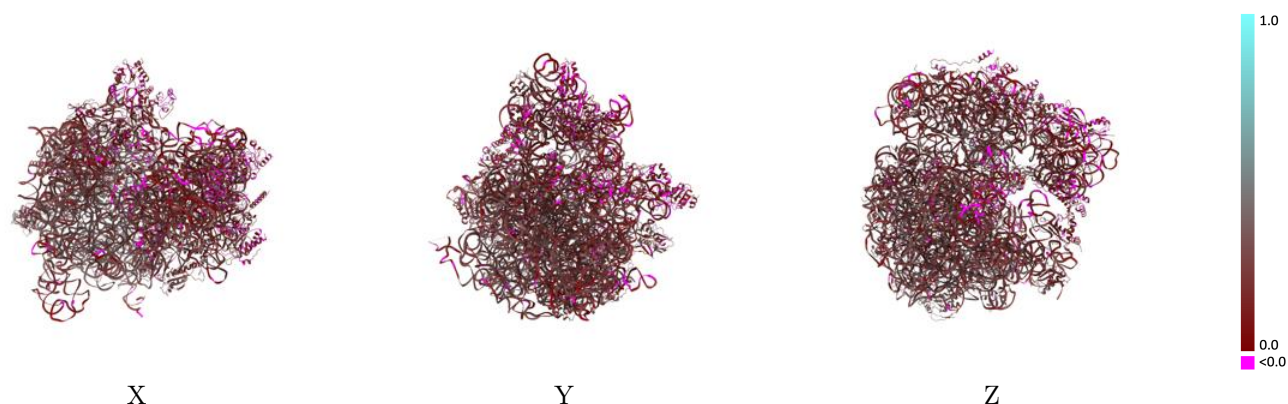
This section contains information regarding the fit between EMDB map EMD-44091 and PDB model 9B1X. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

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



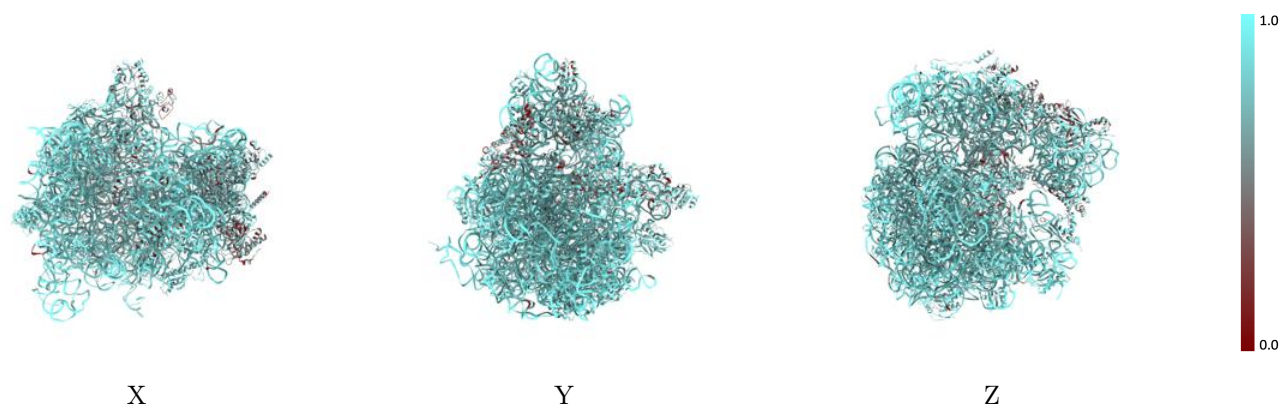
The images above show the 3D surface view of the map at the recommended contour level 0.901 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



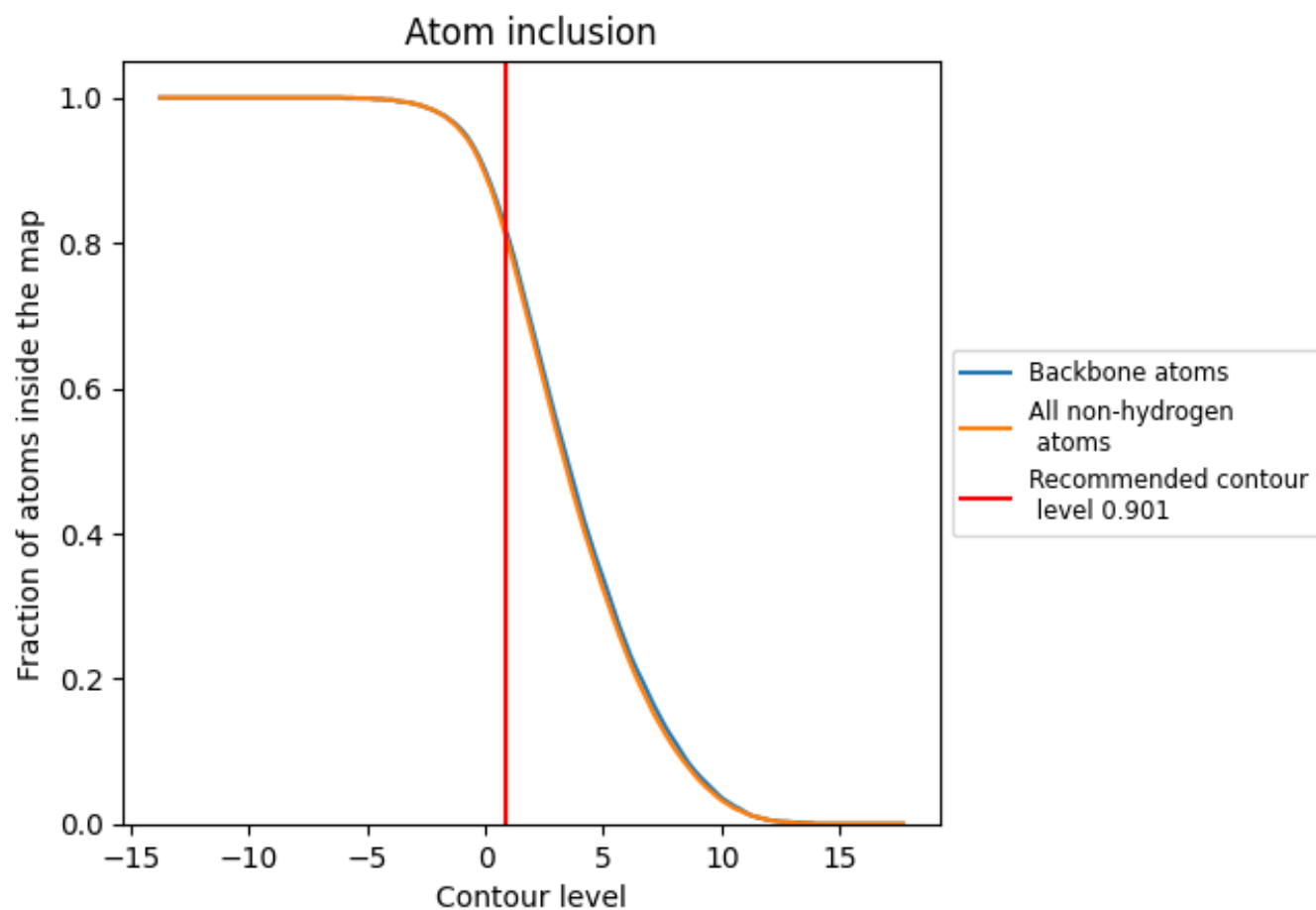
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.901).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 81% 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 (0.901) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8070	 0.2300
1	 0.7880	 0.2900
2	 0.7880	 0.2410
5	 0.6760	 0.2300
A	 0.8410	 0.2020
B	 0.6050	 0.2400
C	 0.5710	 0.1140
D	 0.6330	 0.0950
E	 0.6760	 0.1800
F	 0.7990	 0.2600
G	 0.6580	 0.1710
H	 0.7940	 0.2530
I	 0.7970	 0.2440
J	 0.6050	 0.0900
K	 0.7800	 0.2610
L	 0.7280	 0.2450
M	 0.5820	 0.0750
N	 0.6960	 0.1540
O	 0.8020	 0.2690
P	 0.7610	 0.1910
Q	 0.7850	 0.2290
R	 0.8170	 0.2470
S	 0.7360	 0.1100
T	 0.7890	 0.2060
U	 0.8890	 0.2330
V	 0.4030	 0.1560
X	 0.6840	 0.2000
Y	 0.8500	 0.2540
Z	 0.7510	 0.2600
a	 0.7700	 0.2400
b	 0.8270	 0.3020
c	 0.7520	 0.1560
d	 0.8080	 0.2250
e	 0.5670	 0.0880
f	 0.8080	 0.2670



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Chain	Atom inclusion	Q-score
g	 0.7180	 0.2320
h	 0.8190	 0.2990
i	 0.7210	 0.2210
j	 0.8050	 0.2760
k	 0.8450	 0.2490
l	 0.7650	 0.2340
m	 0.8000	 0.2610
n	 0.8270	 0.2630
o	 0.7960	 0.2770
p	 0.8210	 0.2950
q	 0.8380	 0.2810
r	 0.7350	 0.2010
s	 0.7760	 0.2690
t	 0.7650	 0.2790
u	 0.8100	 0.2710
v	 0.7060	 0.1030
w	 0.7910	 0.2690
x	 0.8280	 0.3010
y	 0.7950	 0.2410
z	 0.8110	 0.3220