



wwPDB EM Validation Summary Report ⓘ

Mar 17, 2026 – 06:31 PM UTC

PDB ID : 6O90 / pdb_00006o90
EMDB ID : EMD-0660
Title : Cryo-EM image reconstruction of the 70S Ribosome *Enterococcus faecalis* Class05
Authors : Jogl, G.; Khayat, R.
Deposited on : 2019-03-12
Resolution : 3.49 Å (reported)
Based on initial models : 4YBB, 5LI0

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

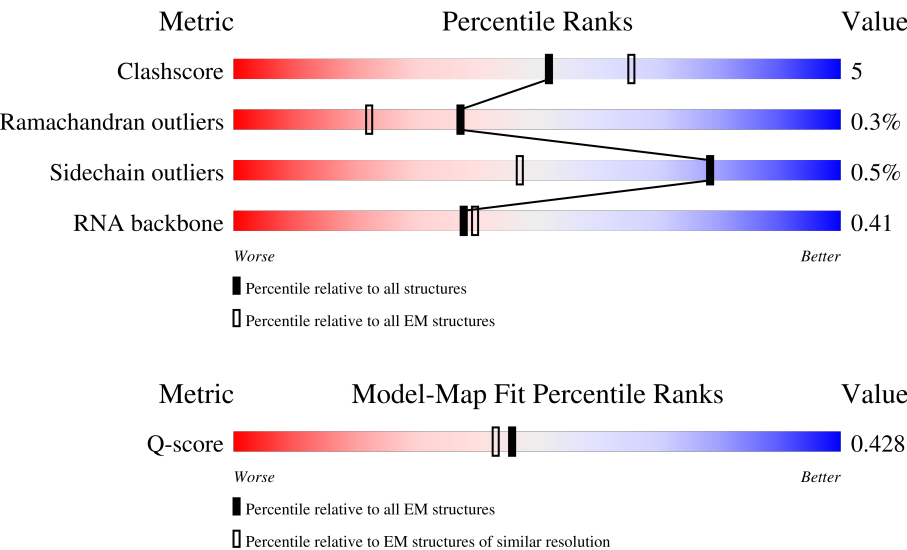
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.49 Å.

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



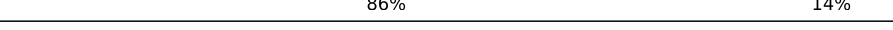
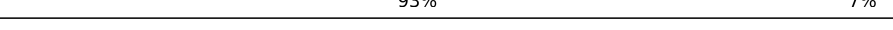
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13600 (2.99 - 3.99)

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

Mol	Chain	Length	Quality of chain
1	a	1528	54% 38% 8%
2	c	204	88% 12%
3	d	201	13% 77% 23%

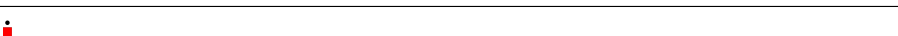
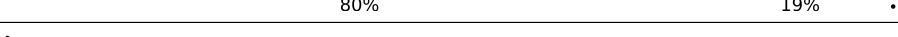
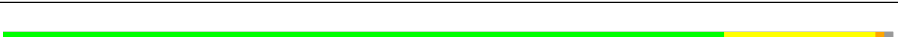
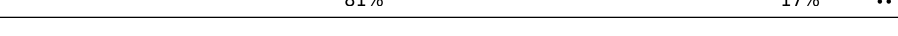
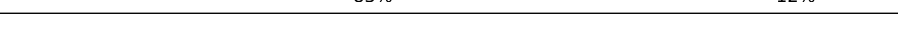
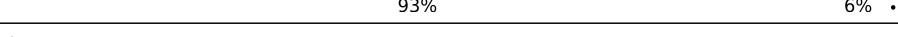
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Mol	Chain	Length	Quality of chain
4	e	163	
5	f	97	
6	g	154	
7	h	131	
8	i	128	
9	j	99	
10	k	117	
11	l	136	
12	m	112	
13	n	60	
14	o	88	
15	p	89	
16	q	83	
17	r	66	
18	s	78	
19	t	81	
20	A	2903	
21	B	116	
22	C	275	
23	D	207	
24	E	206	
25	F	177	
26	G	176	
27	K	145	
28	L	122	

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Mol	Chain	Length	Quality of chain
29	M	146	 85% 13% ..
30	N	141	 6% 80% 19% .
31	O	123	 80% 19% .
32	P	117	 74% 26% .
33	Q	114	 81% 18% .
34	R	118	 81% 17% ..
35	S	102	 84% 16%
36	T	112	 86% 14%
37	U	89	 83% 12% ..
38	V	101	 82% 16% ..
39	W	94	 65% 78% 21% .
40	X	76	 82% 18%
41	Y	54	 93% 6% .
42	Z	61	 90% 7% .
43	0	58	 91% 9%
44	1	83	 80% 20%
45	2	56	 82% 16% .
46	3	49	 86% 14%
47	4	44	 75% 25%
48	5	64	 77% 23%
49	6	38	 74% 26%

2 Entry composition

There are 50 unique types of molecules in this entry. The entry contains 138532 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 RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	a	1528	Total	C	N	O	P	0	0
			32746	14609	5979	10630	1528		

- Molecule 2 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	c	204	Total	C	N	O	S	0	0
			1610	1012	303	292	3		

- Molecule 3 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	d	201	Total	C	N	O	S	0	0
			1620	1016	303	297	4		

- Molecule 4 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	e	163	Total	C	N	O	S	0	0
			1204	759	222	221	2		

- Molecule 5 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	f	97	Total	C	N	O	S	0	0
			795	501	137	154	3		

- Molecule 6 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	g	154	Total	C	N	O	S	0	0
			1229	765	236	222	6		

- Molecule 7 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	h	131	Total	C	N	O	S	0	0
			1041	662	184	193	2		

- Molecule 8 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	i	128	Total	C	N	O	S	0	0
			990	615	197	177	1		

- Molecule 9 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	j	99	Total	C	N	O	S	0	0
			800	504	147	147	2		

- Molecule 10 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	k	117	Total	C	N	O	S	0	0
			863	533	165	161	4		

- Molecule 11 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	l	136	Total	C	N	O	S	0	0
			1065	661	214	188	2		

- Molecule 12 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	m	112	Total	C	N	O	S	0	0
			884	540	180	163	1		

- Molecule 13 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	n	60	Total	C	N	O	S	0	0
			492	310	100	77	5		

- Molecule 14 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	o	88	Total	C	N	O	S	0	0
			741	455	152	133	1		

- Molecule 15 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	p	89	Total	C	N	O	S	0	0
			708	448	131	127	2		

- Molecule 16 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	q	83	Total	C	N	O	S	0	0
			681	427	127	124	3		

- Molecule 17 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	r	66	Total	C	N	O	S	0	0
			537	343	99	94	1		

- Molecule 18 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	s	78	Total	C	N	O	S	0	0
			634	410	113	109	2		

- Molecule 19 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	t	81	Total	C	N	O	S	0	0
			610	372	119	117	2		

- Molecule 20 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	A	2903	Total	C	N	O	P	0	0
			62302	27811	11457	20131	2903		

- Molecule 21 is a RNA chain called 4S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	B	116	Total	C	N	O	P	0	0
			2478	1106	444	812	116		

- Molecule 22 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	C	275	Total	C	N	O	S	0	0
			2114	1310	416	381	7		

- Molecule 23 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	D	207	Total	C	N	O	S	0	0
			1577	992	292	289	4		

- Molecule 24 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	E	206	Total	C	N	O	S	0	0
			1573	984	290	297	2		

- Molecule 25 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	F	177	Total	C	N	O	S	0	0
			1391	887	239	259	6		

- Molecule 26 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	G	176	Total	C	N	O	S	0	0
			1344	842	243	255	4		

- Molecule 27 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	K	145	Total	C	N	O	S	0	0
			1129	713	205	207	4		

- Molecule 28 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	L	122	Total	C	N	O	S	0	0
			922	574	176	170	2		

- Molecule 29 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	M	146	Total	C	N	O	S	0	0
			1094	676	212	205	1		

- Molecule 30 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	N	141	Total	C	N	O	S	0	0
			1117	710	215	185	7		

- Molecule 31 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	O	123	Total	C	N	O	S	0	0
			978	602	190	183	3		

- Molecule 32 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	P	117	Total	C	N	O	S	0	0
			898	556	175	166	1		

- Molecule 33 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms				AltConf	Trace
33	Q	112	Total	C	N	O	0	0
			897	566	177	154		

- Molecule 34 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	R	117	Total	C	N	O	S	0	0
			940	597	181	158	4		

- Molecule 35 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	S	102	Total	C	N	O	S	0	0
			783	499	139	143	2		

- Molecule 36 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	T	112	Total	C	N	O	S	0	0
			849	532	156	159	2		

- Molecule 37 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	U	89	Total	C	N	O	S	0	0
			719	457	127	132	3		

- Molecule 38 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	V	101	Total	C	N	O	S	0	0
			763	486	135	140	2		

- Molecule 39 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	W	94	Total	C	N	O	S	0	0
			757	479	135	139	4		

- Molecule 40 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms				AltConf	Trace
40	X	76	Total	C	N	O	0	0
			571	351	108	112		

- Molecule 41 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	Y	54	Total	C	N	O	S	0	0
			425	265	86	72	2		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	51	ALA	THR	conflict	UNP A0A1B4XRZ8

- Molecule 42 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	Z	61	Total	C	N	O	S	0	0
			504	314	94	95	1		

- Molecule 43 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	0	58	Total	C	N	O	S	0	0
			434	270	81	82	1		

- Molecule 44 is a protein called 50S ribosomal protein L31 type B.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	1	83	Total	C	N	O	S	0	0
			673	424	114	133	2		

- Molecule 45 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	2	56	Total	C	N	O	S	0	0
			429	262	88	73	6		

- Molecule 46 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	3	49	Total	C	N	O	S	0	0
			418	253	85	76	4		

- Molecule 47 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	4	44	Total	C	N	O	S	0	0
			374	227	91	54	2		

- Molecule 48 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	5	64	Total	C	N	O	S	0	0
			522	320	122	78	2		

- Molecule 49 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	6	38	Total	C	N	O	S	0	0
			303	188	66	43	6		

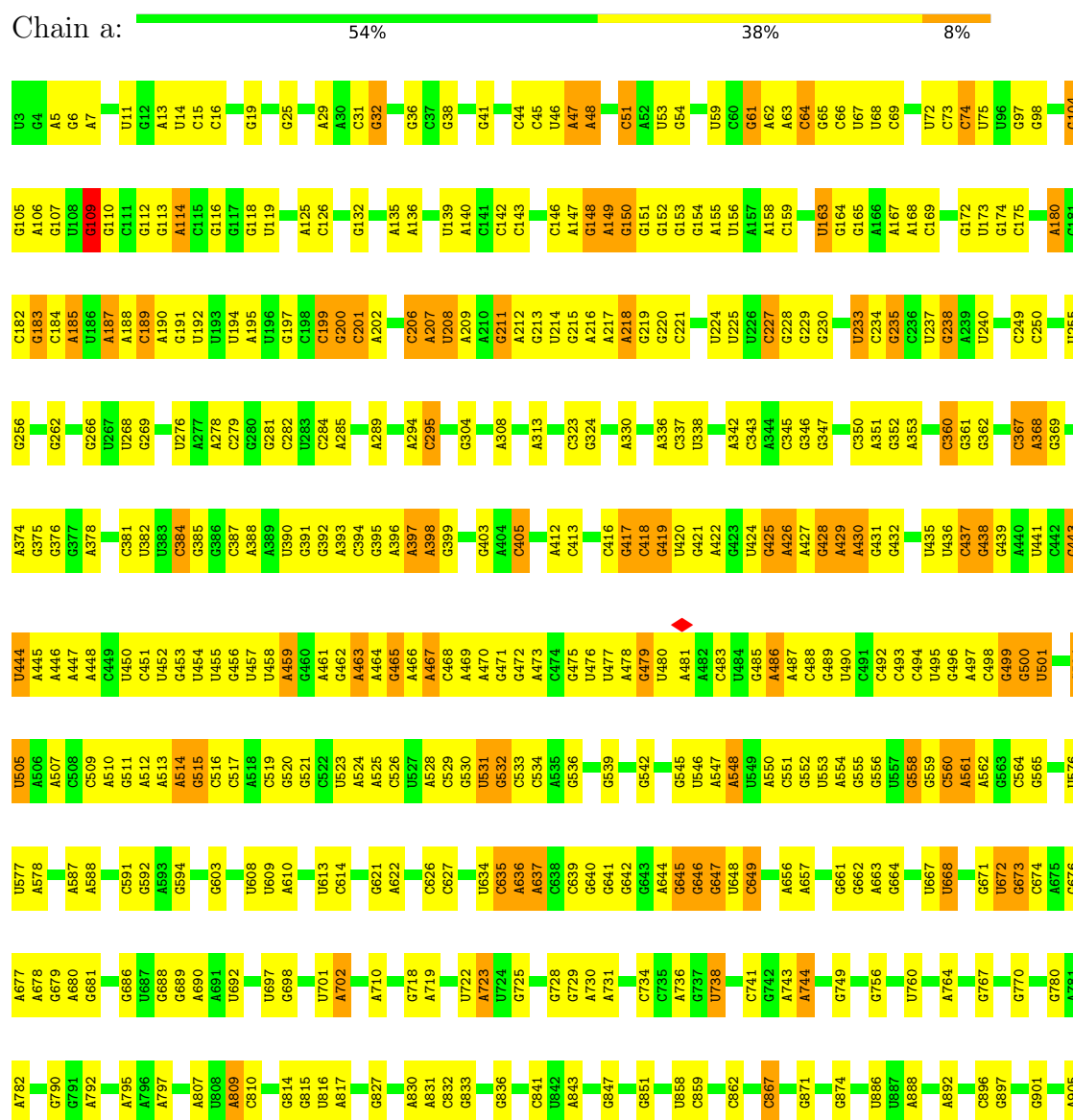
- Molecule 50 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

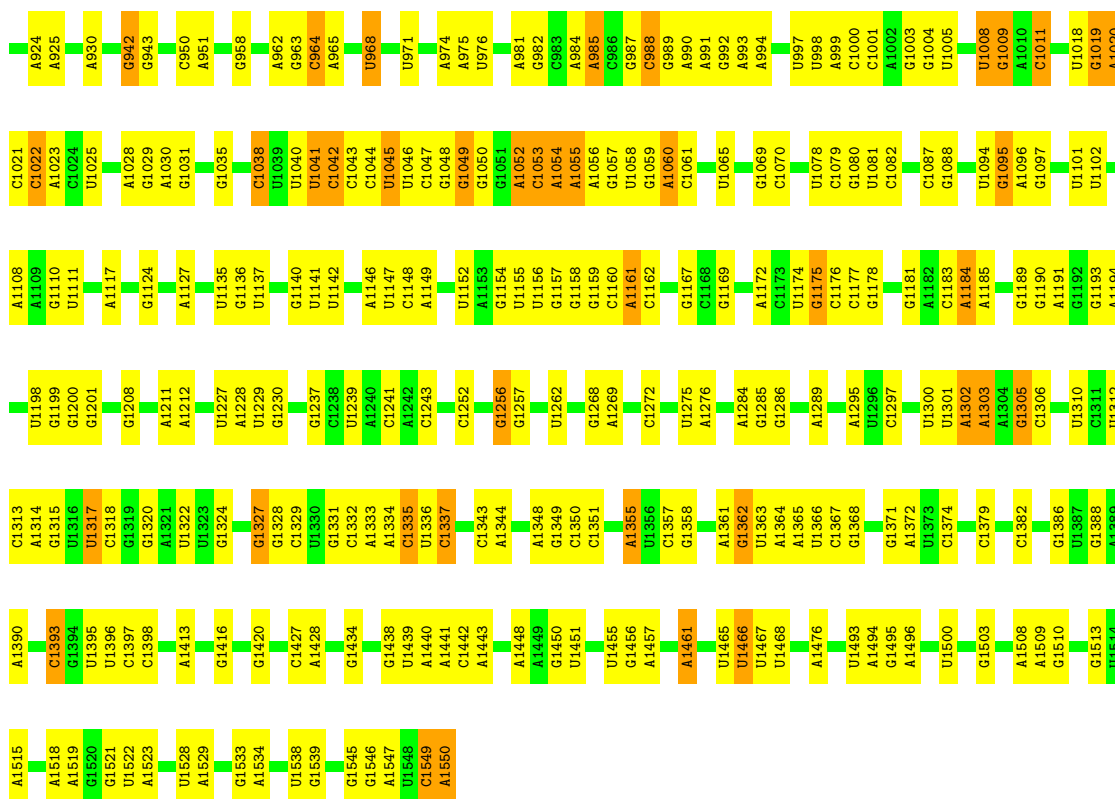
Mol	Chain	Residues	Atoms		AltConf
50	n	1	Total	Zn	0
			1	1	
50	2	1	Total	Zn	0
			1	1	
50	3	1	Total	Zn	0
			1	1	
50	6	1	Total	Zn	0
			1	1	

3 Residue-property plots

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

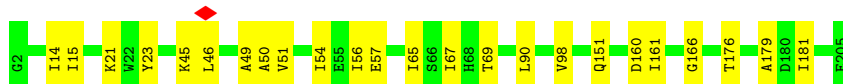
• Molecule 1: 16S rRNA





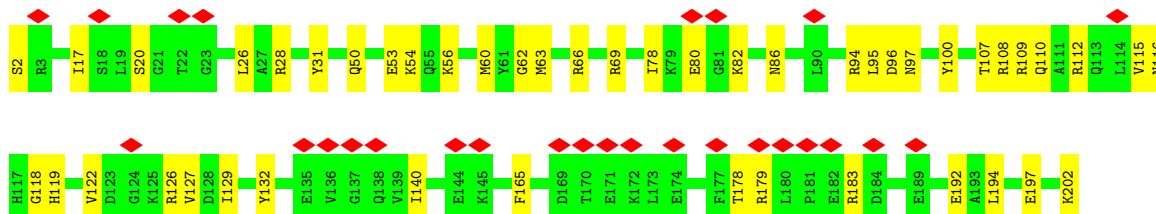
• Molecule 2: 30S ribosomal protein S3

Chain c: 88% 12%



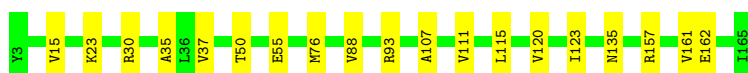
• Molecule 3: 30S ribosomal protein S4

Chain d: 13% 77% 23%



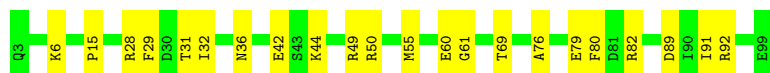
• Molecule 4: 30S ribosomal protein S5

Chain e: 88% 12%



• Molecule 5: 30S ribosomal protein S6

Chain f:  77% 23%



- Molecule 6: 30S ribosomal protein S7

Chain g:  84% 16%



- Molecule 7: 30S ribosomal protein S8

Chain h:  76% 24%



- Molecule 8: 30S ribosomal protein S9

Chain i:  75% 23%



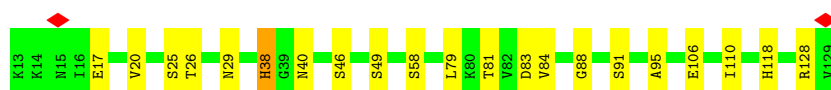
- Molecule 9: 30S ribosomal protein S10

Chain j:  76% 24%



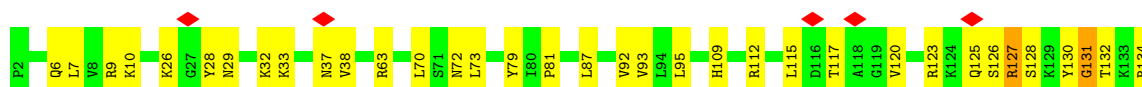
- Molecule 10: 30S ribosomal protein S11

Chain k:  82% 17%



- Molecule 11: 30S ribosomal protein S12

Chain l:  5% 74% 24%





- Molecule 12: 30S ribosomal protein S13

Chain m: 86% 14%



- Molecule 13: 30S ribosomal protein S14 type Z

Chain n: 90% 10%



- Molecule 14: 30S ribosomal protein S15

Chain o: 86% 14%



- Molecule 15: 30S ribosomal protein S16

Chain p: 76% 24%



- Molecule 16: 30S ribosomal protein S17

Chain q: 83% 17%



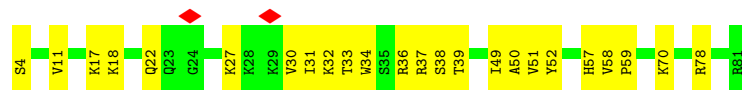
- Molecule 17: 30S ribosomal protein S18

Chain r: 86% 14%

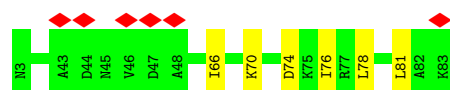


- Molecule 18: 30S ribosomal protein S19

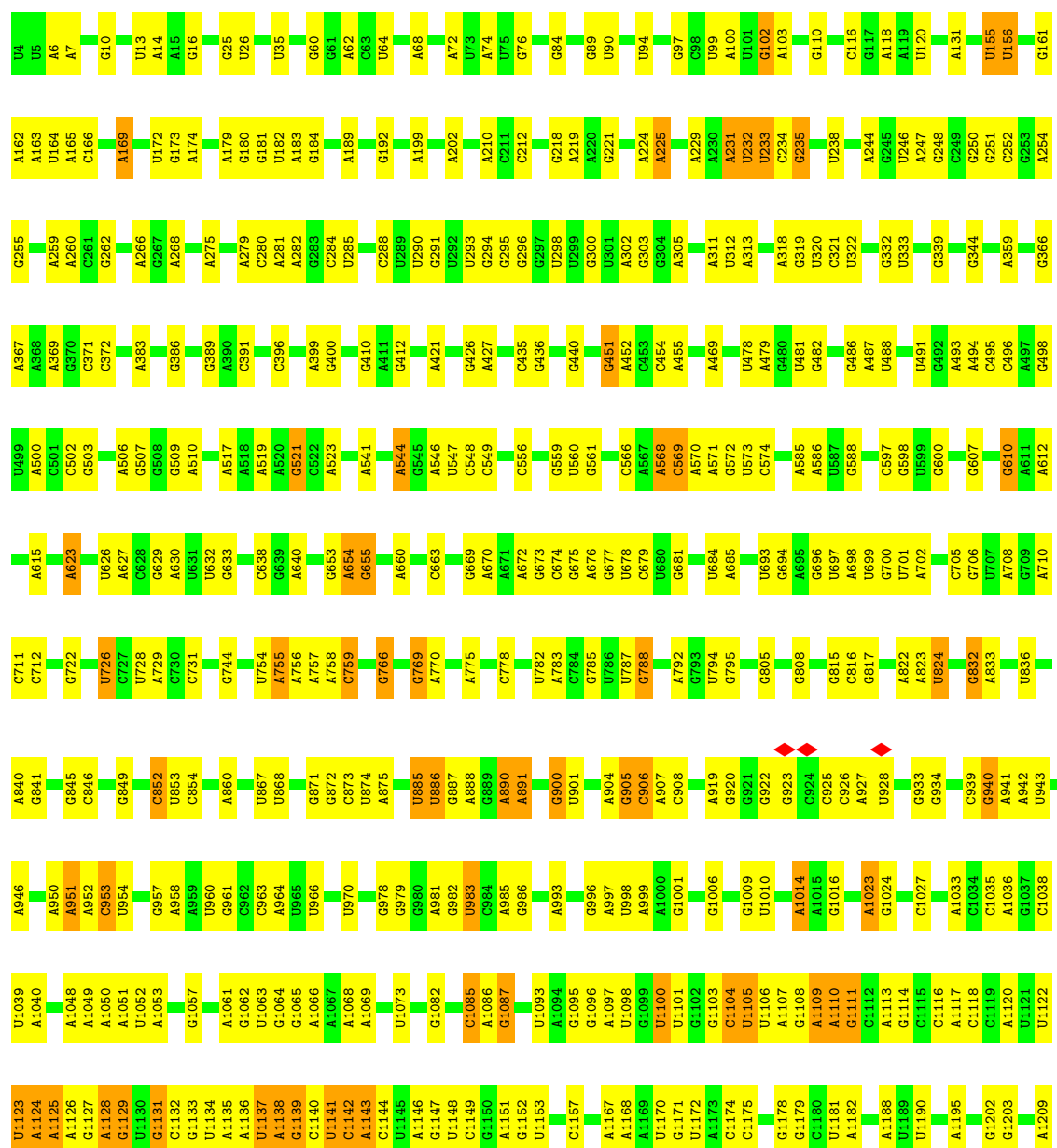
Chain s: 69% 31%

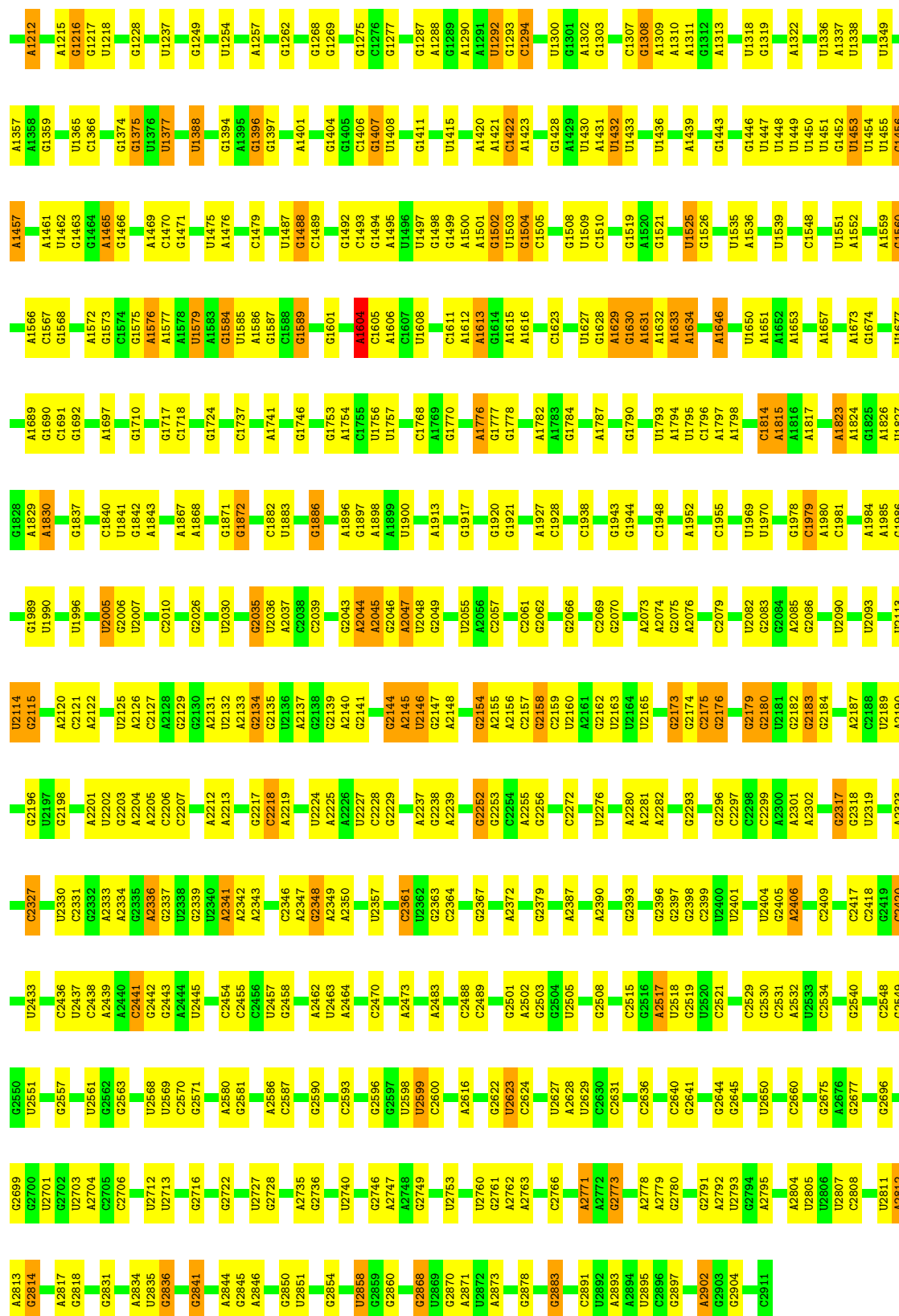


- Molecule 19: 30S ribosomal protein S20



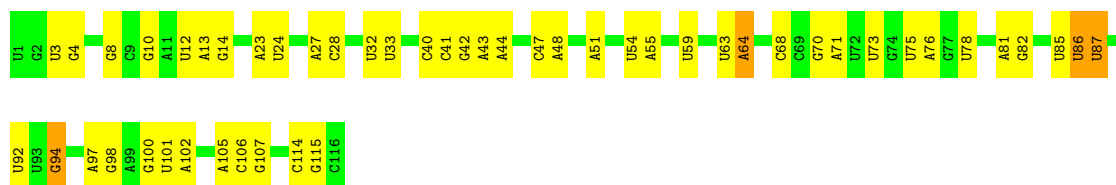
- Molecule 20: 23S rRNA





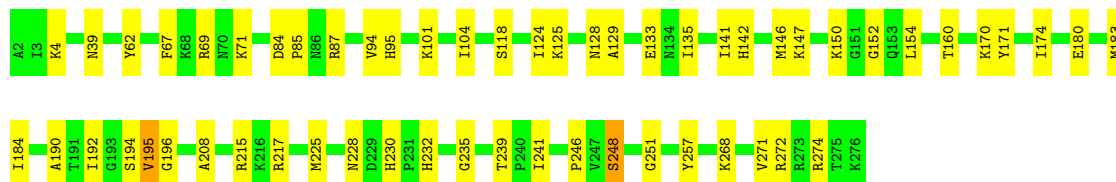
- Molecule 21: 4S rRNA





- Molecule 22: 50S ribosomal protein L2

Chain C: 79% 20% .



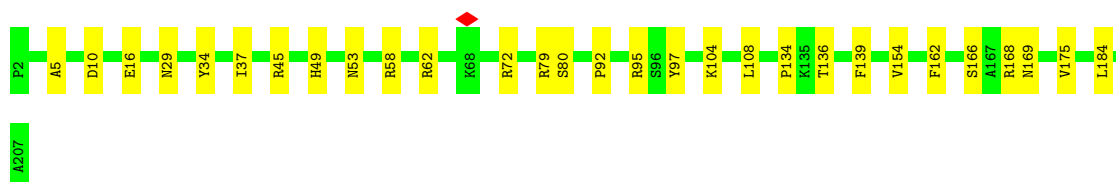
- Molecule 23: 50S ribosomal protein L3

Chain D: 84% 15% .



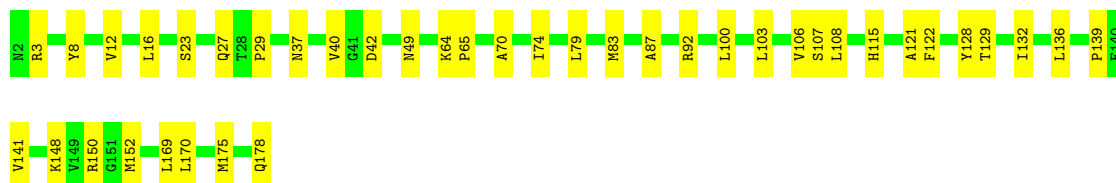
- Molecule 24: 50S ribosomal protein L4

Chain E: 86% 14%



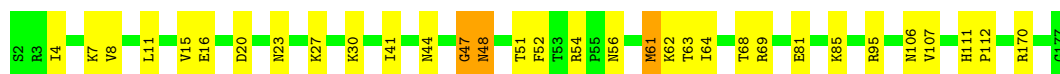
- Molecule 25: 50S ribosomal protein L5

Chain F: 77% 23%



- Molecule 26: 50S ribosomal protein L6

Chain G: 82% 16% .



- Molecule 27: 50S ribosomal protein L13

Chain K:  88% 10%



- Molecule 28: 50S ribosomal protein L14

Chain L:  91% 9%



- Molecule 29: 50S ribosomal protein L15

Chain M:  85% 13%



- Molecule 30: 50S ribosomal protein L16

Chain N:  6% 80% 19%



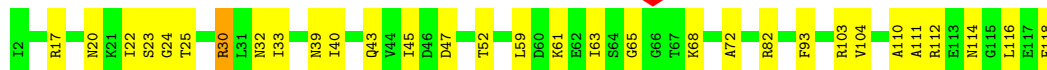
- Molecule 31: 50S ribosomal protein L17

Chain O:  80% 19%



- Molecule 32: 50S ribosomal protein L18

Chain P:  74% 26%



- Molecule 33: 50S ribosomal protein L19

Chain Q:  81% 18%



- Molecule 34: 50S ribosomal protein L20

Chain R:  81% 17% ..



- Molecule 35: 50S ribosomal protein L21

Chain S:  84% 16%



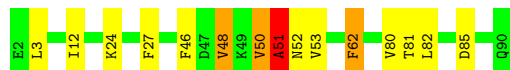
- Molecule 36: 50S ribosomal protein L22

Chain T:  86% 14%



- Molecule 37: 50S ribosomal protein L23

Chain U:  83% 12% ..



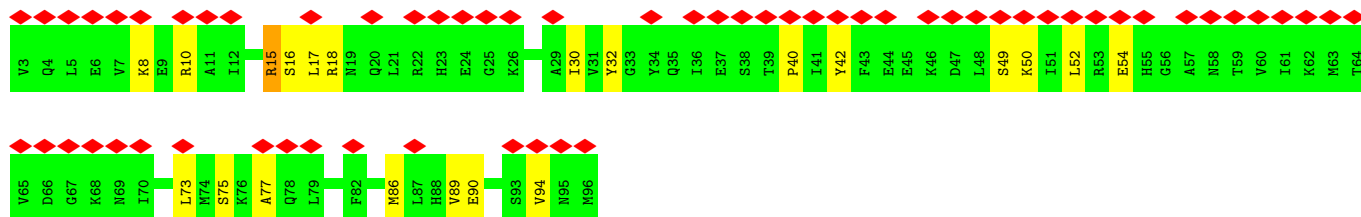
- Molecule 38: 50S ribosomal protein L24

Chain V:  82% 16% ..



- Molecule 39: 50S ribosomal protein L25

Chain W:  65% 78% 21% .



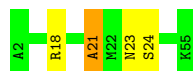
- Molecule 40: 50S ribosomal protein L27

Chain X:  82% 18%



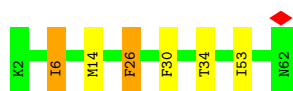
- Molecule 41: 50S ribosomal protein L28

Chain Y: 93% 6% .



- Molecule 42: 50S ribosomal protein L29

Chain Z: 90% 7% .



- Molecule 43: 50S ribosomal protein L30

Chain 0: 91% 9%



- Molecule 44: 50S ribosomal protein L31 type B

Chain 1: 80% 20%



- Molecule 45: 50S ribosomal protein L32

Chain 2: 82% 16% .



- Molecule 46: 50S ribosomal protein L33

Chain 3: 86% 14%

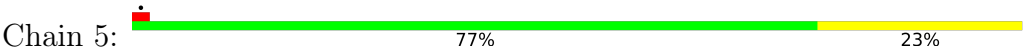


- Molecule 47: 50S ribosomal protein L34

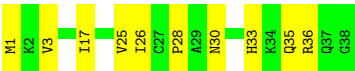
Chain 4: 75% 25%



• Molecule 48: 50S ribosomal protein L35



• Molecule 49: 50S ribosomal protein L36



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	46244	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 ($e^-/\text{\AA}^2$)	25	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	2.323	Depositor
Minimum map value	-0.893	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.095	Depositor
Recommended contour level	0.238	Depositor
Map size (Å)	482.68, 482.68, 482.68	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.097, 1.097, 1.097	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.31	0/36657	0.45	1/57173 (0.0%)
2	c	0.24	0/1635	0.54	0/2197
3	d	0.29	0/1650	0.71	0/2217
4	e	0.29	0/1217	0.62	2/1641 (0.1%)
5	f	0.23	0/807	0.50	0/1087
6	g	0.23	0/1249	0.55	0/1682
7	h	0.29	0/1054	0.56	0/1417
8	i	0.26	0/1003	0.64	2/1343 (0.1%)
9	j	0.28	0/812	0.66	0/1093
10	k	0.26	0/878	0.57	2/1185 (0.2%)
11	l	0.35	0/1082	0.81	1/1453 (0.1%)
12	m	0.23	0/890	0.64	0/1195
13	n	0.26	0/504	0.56	0/669
14	o	0.29	0/751	0.57	0/1001
15	p	0.34	0/720	0.65	0/966
16	q	0.32	0/689	0.65	0/920
17	r	0.28	0/544	0.61	0/728
18	s	0.27	0/650	0.64	0/872
19	t	0.31	0/612	0.63	0/818
20	A	0.41	0/69785	0.47	1/108842 (0.0%)
21	B	0.33	0/2770	0.44	0/4311
22	C	0.49	1/2148 (0.0%)	0.77	2/2888 (0.1%)
23	D	0.46	0/1597	0.75	0/2143
24	E	0.42	0/1595	0.65	0/2157
25	F	0.29	0/1410	0.58	0/1895
26	G	0.33	0/1362	0.69	1/1831 (0.1%)
27	K	0.48	0/1148	0.71	0/1546
28	L	0.48	0/929	0.68	0/1247
29	M	0.40	0/1102	0.91	3/1467 (0.2%)
30	N	0.43	0/1139	0.81	2/1515 (0.1%)
31	O	0.48	0/984	0.85	2/1317 (0.2%)
32	P	0.38	0/907	0.67	0/1214

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	Q	0.45	0/911	0.67	0/1227
34	R	0.50	0/951	0.73	1/1260 (0.1%)
35	S	0.42	0/794	0.68	0/1064
36	T	0.44	0/858	0.67	0/1157
37	U	0.41	0/725	0.81	2/969 (0.2%)
38	V	0.35	0/772	0.82	2/1035 (0.2%)
39	W	0.25	0/768	0.67	0/1032
40	X	0.47	0/576	0.71	0/768
41	Y	0.31	0/431	0.62	0/574
42	Z	0.36	0/505	0.61	0/672
43	0	0.42	0/434	0.68	0/583
44	1	0.26	0/690	0.77	0/930
45	2	0.48	0/436	0.70	1/578 (0.2%)
46	3	0.30	0/422	0.53	0/561
47	4	0.49	0/377	0.70	0/491
48	5	0.42	0/528	0.70	0/689
49	6	0.44	0/308	0.66	0/407
All	All	0.38	1/150766 (0.0%)	0.53	25/226027 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
11	l	0	2
12	m	0	1
16	q	0	1
19	t	0	1
22	C	0	2
23	D	0	2
24	E	0	2
25	F	0	1
26	G	0	2
27	K	0	2
29	M	0	3
30	N	0	2
31	O	0	2
32	P	0	3
35	S	0	2
37	U	0	2
38	V	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
39	W	0	1
40	X	0	1
41	Y	0	1
47	4	0	1
All	All	0	36

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	C	248	SER	C-N	5.19	1.38	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
38	V	71	ILE	CA-C-N	6.92	138.68	121.80
38	V	71	ILE	C-N-CA	6.92	138.68	121.80
26	G	4	ILE	N-CA-C	-6.77	106.19	112.96
31	O	41	ARG	CA-CB-CG	6.24	126.59	114.10
10	k	38	HIS	CA-C-N	6.07	132.62	121.70

There are no chirality outliers.

5 of 36 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
11	l	130	TYR	Peptide
11	l	37	ASN	Peptide
12	m	43	VAL	Peptide
16	q	73	THR	Peptide
19	t	66	ILE	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	a	32746	0	16491	336	0
2	c	1610	0	1653	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	d	1620	0	1641	32	0
4	e	1204	0	1280	11	0
5	f	795	0	789	14	0
6	g	1229	0	1259	18	0
7	h	1041	0	1092	19	0
8	i	990	0	1035	22	0
9	j	800	0	846	15	0
10	k	863	0	882	12	0
11	l	1065	0	1133	23	0
12	m	884	0	932	11	0
13	n	492	0	510	5	0
14	o	741	0	756	9	0
15	p	708	0	746	19	0
16	q	681	0	719	10	0
17	r	537	0	572	5	0
18	s	634	0	649	17	0
19	t	610	0	640	3	0
20	A	62302	0	31302	356	0
21	B	2478	0	1248	21	0
22	C	2114	0	2197	38	0
23	D	1577	0	1647	24	0
24	E	1573	0	1621	19	0
25	F	1391	0	1454	24	0
26	G	1344	0	1372	22	0
27	K	1129	0	1163	14	0
28	L	922	0	985	7	0
29	M	1094	0	1145	14	0
30	N	1117	0	1177	18	0
31	O	978	0	1017	17	0
32	P	898	0	932	18	0
33	Q	897	0	943	15	0
34	R	940	0	998	15	0
35	S	783	0	821	10	0
36	T	849	0	906	10	0
37	U	719	0	769	8	0
38	V	763	0	824	9	0
39	W	757	0	780	13	0
40	X	571	0	578	8	0
41	Y	425	0	460	3	0
42	Z	504	0	538	5	0
43	0	434	0	462	8	0
44	1	673	0	639	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
45	2	429	0	446	8	0
46	3	418	0	433	5	0
47	4	374	0	424	9	0
48	5	522	0	576	19	0
49	6	303	0	338	6	0
50	2	1	0	0	0	0
50	3	1	0	0	0	0
50	6	1	0	0	0	0
50	n	1	0	0	0	0
All	All	138532	0	91820	1166	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 1166 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:A:2893:A:C2	20:A:2893:A:C4	2.20	1.18
20:A:1100:U:N3	20:A:1128:A:C8	2.19	1.10
43:0:50:ILE:CG2	43:0:50:ILE:CG1	2.30	1.10
1:a:150:G:H21	1:a:185:A:N6	1.49	1.07
1:a:154:G:H1	1:a:180:A:N6	1.55	1.04

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	c	202/204 (99%)	178 (88%)	24 (12%)	0	100	100
3	d	199/201 (99%)	169 (85%)	30 (15%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	e	161/163 (99%)	142 (88%)	19 (12%)	0	100	100
5	f	95/97 (98%)	88 (93%)	7 (7%)	0	100	100
6	g	152/154 (99%)	138 (91%)	14 (9%)	0	100	100
7	h	129/131 (98%)	116 (90%)	13 (10%)	0	100	100
8	i	126/128 (98%)	110 (87%)	15 (12%)	1 (1%)	16	49
9	j	97/99 (98%)	87 (90%)	10 (10%)	0	100	100
10	k	115/117 (98%)	98 (85%)	17 (15%)	0	100	100
11	l	134/136 (98%)	106 (79%)	27 (20%)	1 (1%)	18	51
12	m	110/112 (98%)	86 (78%)	24 (22%)	0	100	100
13	n	58/60 (97%)	53 (91%)	5 (9%)	0	100	100
14	o	86/88 (98%)	79 (92%)	7 (8%)	0	100	100
15	p	87/89 (98%)	75 (86%)	12 (14%)	0	100	100
16	q	81/83 (98%)	67 (83%)	14 (17%)	0	100	100
17	r	64/66 (97%)	53 (83%)	11 (17%)	0	100	100
18	s	76/78 (97%)	60 (79%)	16 (21%)	0	100	100
19	t	79/81 (98%)	71 (90%)	8 (10%)	0	100	100
22	C	273/275 (99%)	236 (86%)	37 (14%)	0	100	100
23	D	205/207 (99%)	182 (89%)	22 (11%)	1 (0%)	24	57
24	E	204/206 (99%)	175 (86%)	29 (14%)	0	100	100
25	F	175/177 (99%)	149 (85%)	26 (15%)	0	100	100
26	G	172/176 (98%)	145 (84%)	25 (14%)	2 (1%)	10	41
27	K	142/145 (98%)	124 (87%)	18 (13%)	0	100	100
28	L	120/122 (98%)	98 (82%)	22 (18%)	0	100	100
29	M	143/146 (98%)	105 (73%)	35 (24%)	3 (2%)	5	31
30	N	137/141 (97%)	111 (81%)	26 (19%)	0	100	100
31	O	119/123 (97%)	95 (80%)	24 (20%)	0	100	100
32	P	115/117 (98%)	100 (87%)	15 (13%)	0	100	100
33	Q	110/114 (96%)	98 (89%)	12 (11%)	0	100	100
34	R	112/118 (95%)	108 (96%)	4 (4%)	0	100	100
35	S	100/102 (98%)	92 (92%)	8 (8%)	0	100	100
36	T	110/112 (98%)	95 (86%)	15 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
37	U	87/89 (98%)	72 (83%)	13 (15%)	2 (2%)	5	30
38	V	99/101 (98%)	75 (76%)	21 (21%)	3 (3%)	3	25
39	W	92/94 (98%)	74 (80%)	17 (18%)	1 (1%)	11	43
40	X	73/76 (96%)	64 (88%)	9 (12%)	0	100	100
41	Y	52/54 (96%)	44 (85%)	8 (15%)	0	100	100
42	Z	59/61 (97%)	55 (93%)	4 (7%)	0	100	100
43	0	56/58 (97%)	53 (95%)	3 (5%)	0	100	100
44	1	81/83 (98%)	60 (74%)	21 (26%)	0	100	100
45	2	54/56 (96%)	47 (87%)	7 (13%)	0	100	100
46	3	47/49 (96%)	45 (96%)	2 (4%)	0	100	100
47	4	42/44 (96%)	40 (95%)	2 (5%)	0	100	100
48	5	62/64 (97%)	55 (89%)	7 (11%)	0	100	100
49	6	36/38 (95%)	33 (92%)	3 (8%)	0	100	100
All	All	5128/5235 (98%)	4406 (86%)	708 (14%)	14 (0%)	37	67

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
26	G	48	ASN
29	M	95	VAL
29	M	98	GLU
39	W	16	SER
11	l	131	GLY

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	c	162/162 (100%)	162 (100%)	0	100	100
3	d	175/175 (100%)	174 (99%)	1 (1%)	78	79
4	e	126/126 (100%)	126 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	f	86/86 (100%)	86 (100%)	0	100	100
6	g	131/131 (100%)	131 (100%)	0	100	100
7	h	112/112 (100%)	112 (100%)	0	100	100
8	i	101/101 (100%)	100 (99%)	1 (1%)	68	75
9	j	90/90 (100%)	90 (100%)	0	100	100
10	k	91/91 (100%)	91 (100%)	0	100	100
11	l	118/118 (100%)	117 (99%)	1 (1%)	73	77
12	m	95/95 (100%)	95 (100%)	0	100	100
13	n	51/51 (100%)	50 (98%)	1 (2%)	48	67
14	o	78/78 (100%)	78 (100%)	0	100	100
15	p	79/79 (100%)	79 (100%)	0	100	100
16	q	76/76 (100%)	76 (100%)	0	100	100
17	r	57/57 (100%)	57 (100%)	0	100	100
18	s	68/68 (100%)	67 (98%)	1 (2%)	57	71
19	t	62/62 (100%)	61 (98%)	1 (2%)	55	70
22	C	224/225 (100%)	223 (100%)	1 (0%)	84	81
23	D	168/170 (99%)	167 (99%)	1 (1%)	78	79
24	E	172/172 (100%)	171 (99%)	1 (1%)	78	79
25	F	154/154 (100%)	154 (100%)	0	100	100
26	G	145/146 (99%)	144 (99%)	1 (1%)	76	78
27	K	121/122 (99%)	120 (99%)	1 (1%)	73	77
28	L	98/98 (100%)	97 (99%)	1 (1%)	68	75
29	M	111/112 (99%)	110 (99%)	1 (1%)	70	76
30	N	112/112 (100%)	112 (100%)	0	100	100
31	O	105/105 (100%)	105 (100%)	0	100	100
32	P	91/91 (100%)	91 (100%)	0	100	100
33	Q	94/97 (97%)	94 (100%)	0	100	100
34	R	92/94 (98%)	92 (100%)	0	100	100
35	S	82/83 (99%)	82 (100%)	0	100	100
36	T	95/95 (100%)	95 (100%)	0	100	100
37	U	80/80 (100%)	76 (95%)	4 (5%)	22	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
38	V	85/85 (100%)	85 (100%)	0	100	100
39	W	85/85 (100%)	85 (100%)	0	100	100
40	X	60/61 (98%)	60 (100%)	0	100	100
41	Y	47/47 (100%)	47 (100%)	0	100	100
42	Z	55/55 (100%)	53 (96%)	2 (4%)	31	57
43	0	48/49 (98%)	48 (100%)	0	100	100
44	1	75/75 (100%)	75 (100%)	0	100	100
45	2	46/46 (100%)	45 (98%)	1 (2%)	45	65
46	3	49/49 (100%)	49 (100%)	0	100	100
47	4	39/39 (100%)	39 (100%)	0	100	100
48	5	51/51 (100%)	51 (100%)	0	100	100
49	6	35/35 (100%)	35 (100%)	0	100	100
All	All	4377/4391 (100%)	4357 (100%)	20 (0%)	78	80

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
37	U	48	VAL
42	Z	6	ILE
45	2	24	ILE
42	Z	26	PHE
22	C	257	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 74 such sidechains are listed below:

Mol	Chain	Res	Type
40	X	87	GLN
48	5	34	HIS
41	Y	17	ASN
44	1	41	ASN
18	s	22	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	a	1526/1528 (99%)	441 (28%)	0

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
20	A	2895/2903 (99%)	650 (22%)	20 (0%)
21	B	113/116 (97%)	26 (23%)	2 (1%)
All	All	4534/4547 (99%)	1117 (24%)	22 (0%)

5 of 1117 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	a	5	A
1	a	6	G
1	a	7	A
1	a	11	U
1	a	13	A

5 of 22 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
20	A	1604	A
20	A	2281	A
20	A	1628	G
20	A	2317	G
20	A	654	A

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 4 ligands modelled in this entry, 4 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
20	A	2
1	a	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	928:U	O3'	931:C	P	16.19
1	a	75:U	O3'	96:U	P	13.98
1	A	1579:U	O3'	1583:A	P	9.00

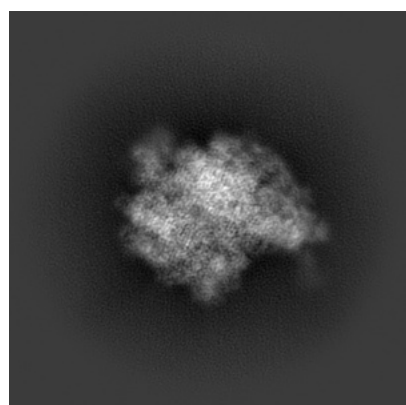
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-0660. These allow visual inspection of the internal detail of the map and identification of artifacts.

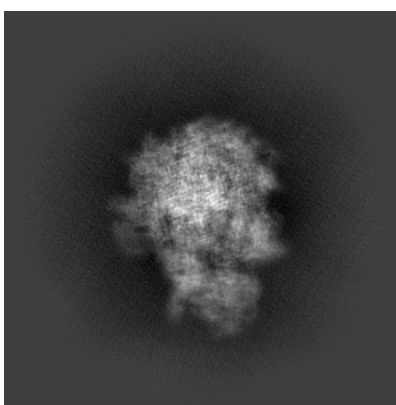
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

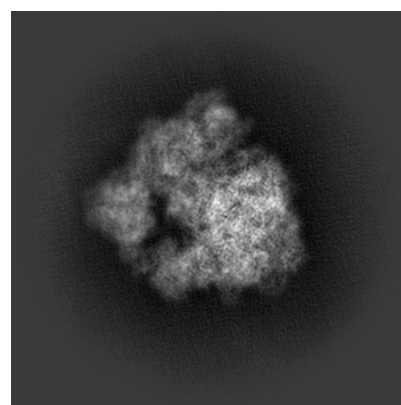
6.1.1 Primary map



X



Y

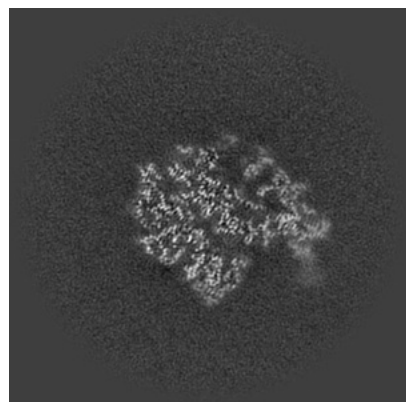


Z

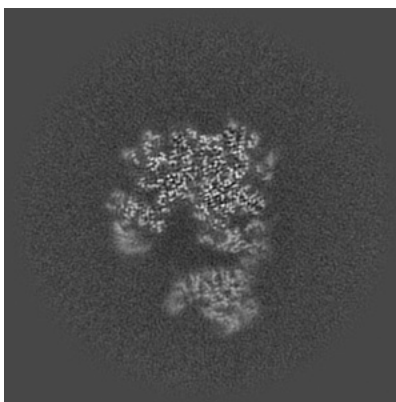
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

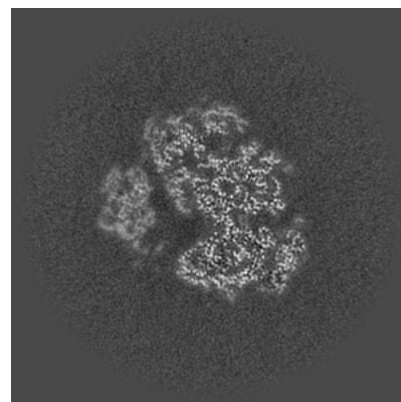
6.2.1 Primary map



X Index: 220



Y Index: 220

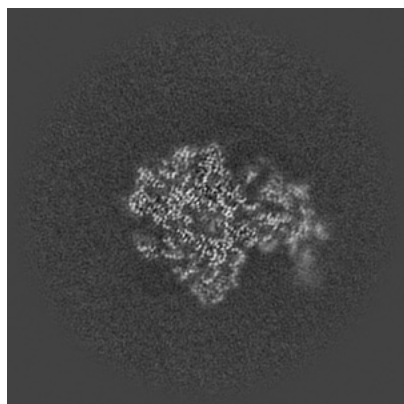


Z Index: 220

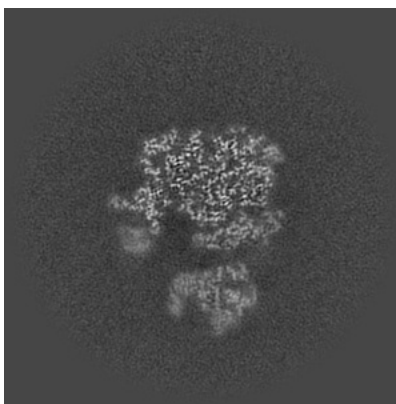
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

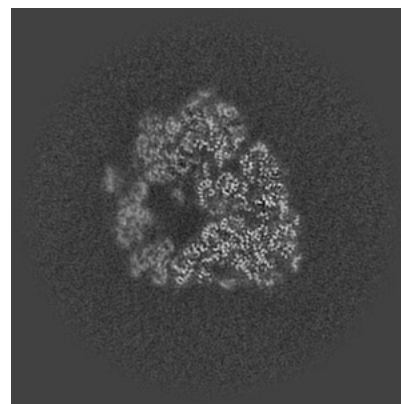
6.3.1 Primary map



X Index: 226



Y Index: 231

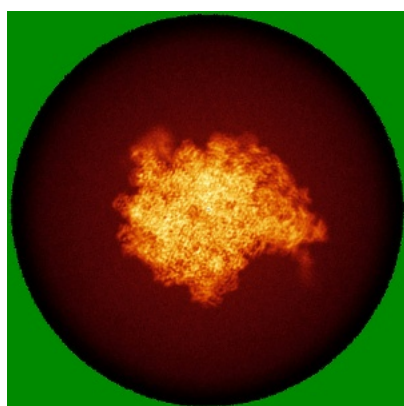


Z Index: 209

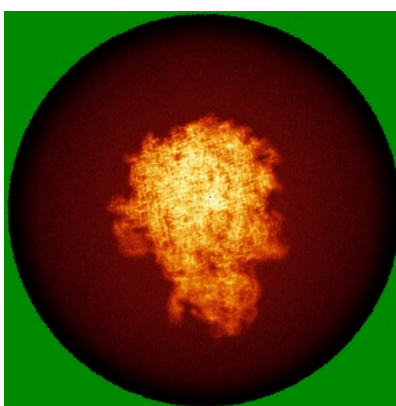
The images above show the largest variance slices of the map in three orthogonal directions.

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

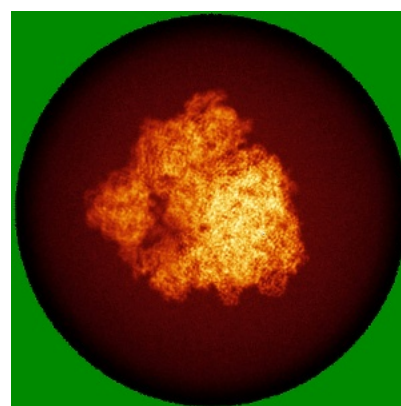
6.4.1 Primary map



X



Y

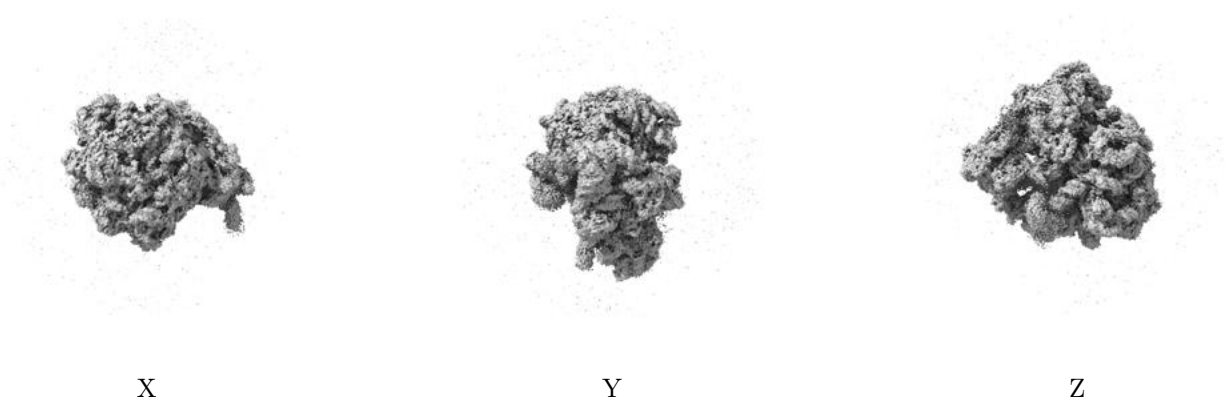


Z

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 0.238. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

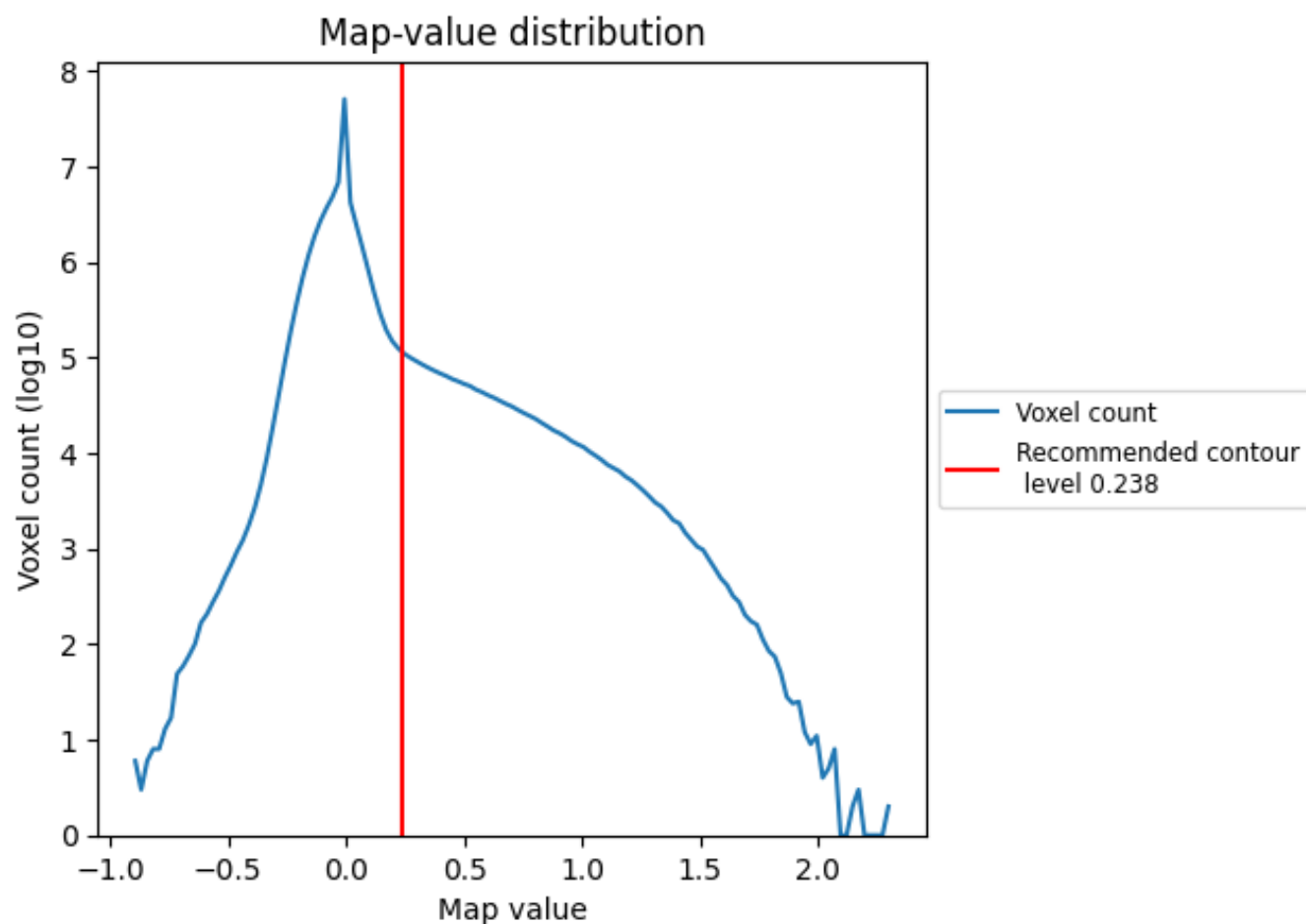
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

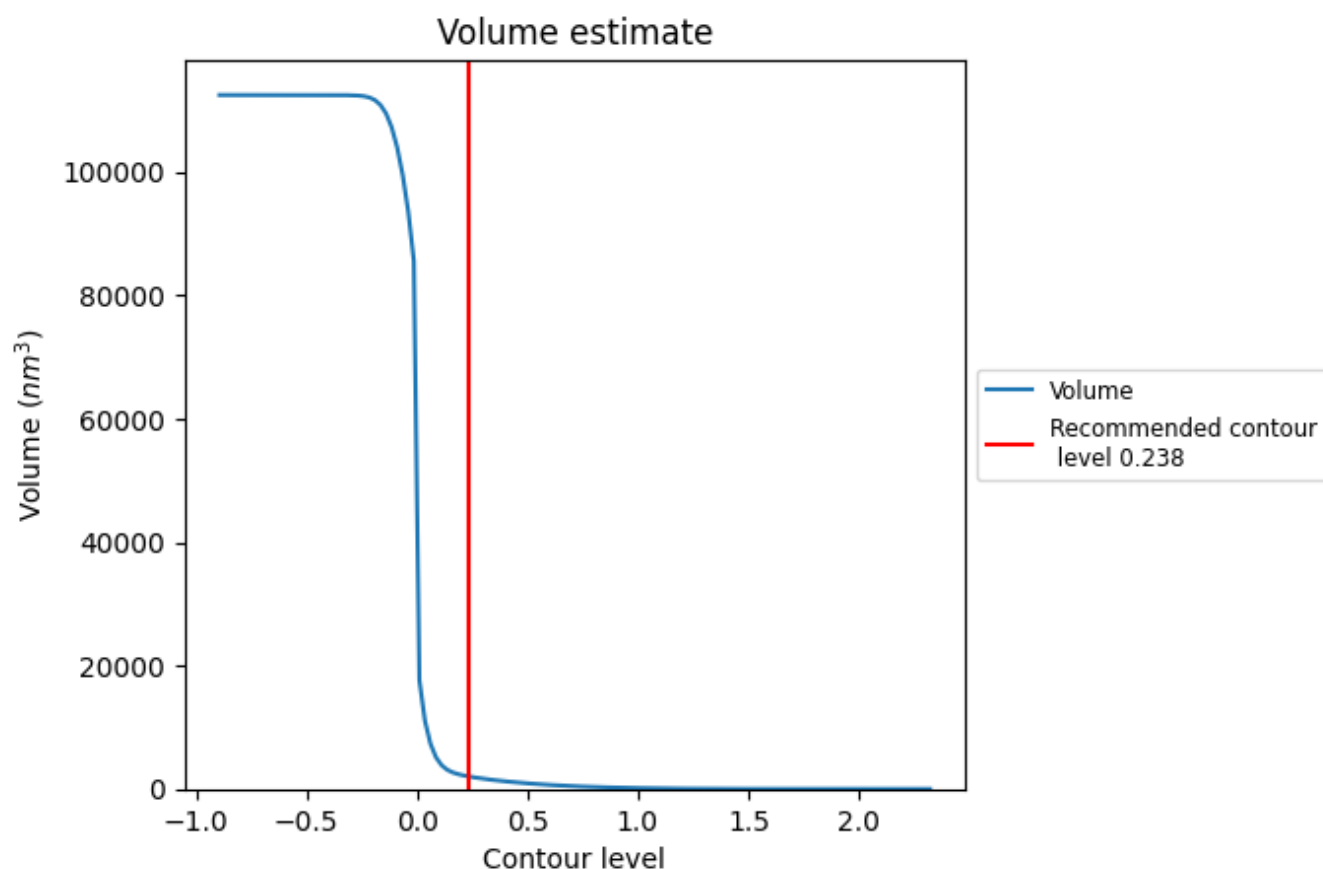
This section contains the results of statistical analysis of the map.

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.

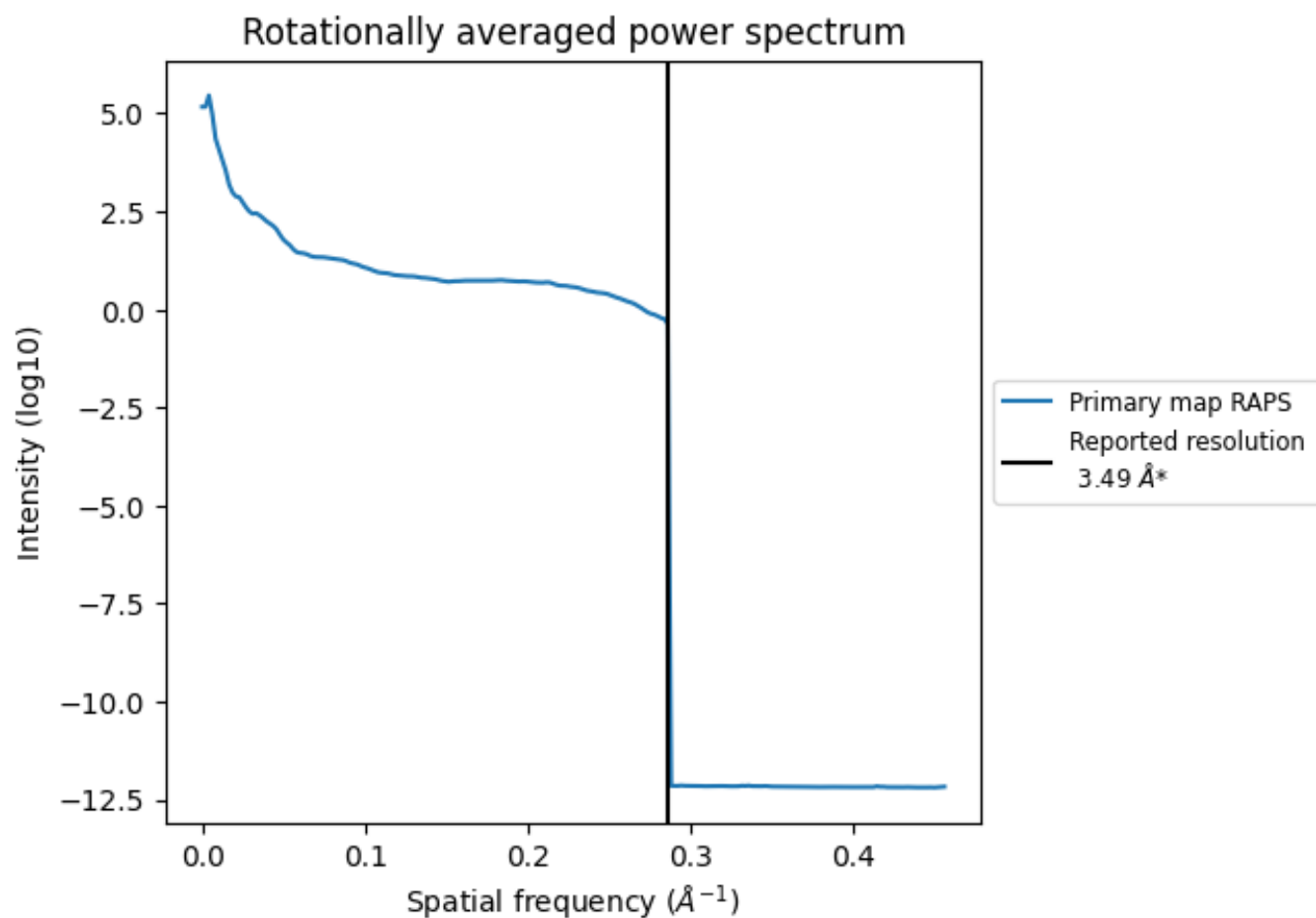
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2006 nm³; this corresponds to an approximate mass of 1813 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 0.287 $\text{\AA}^{-1}$

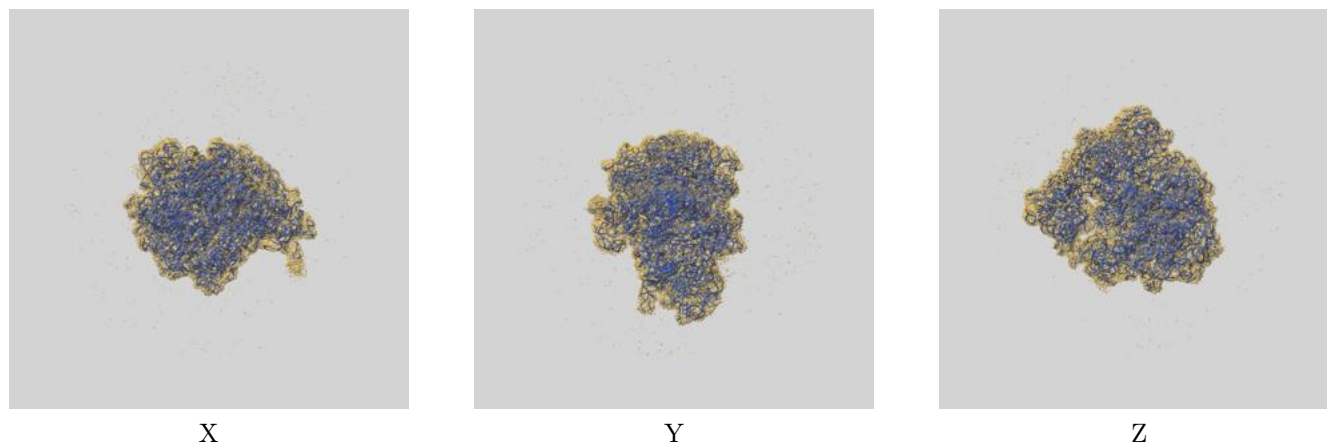
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

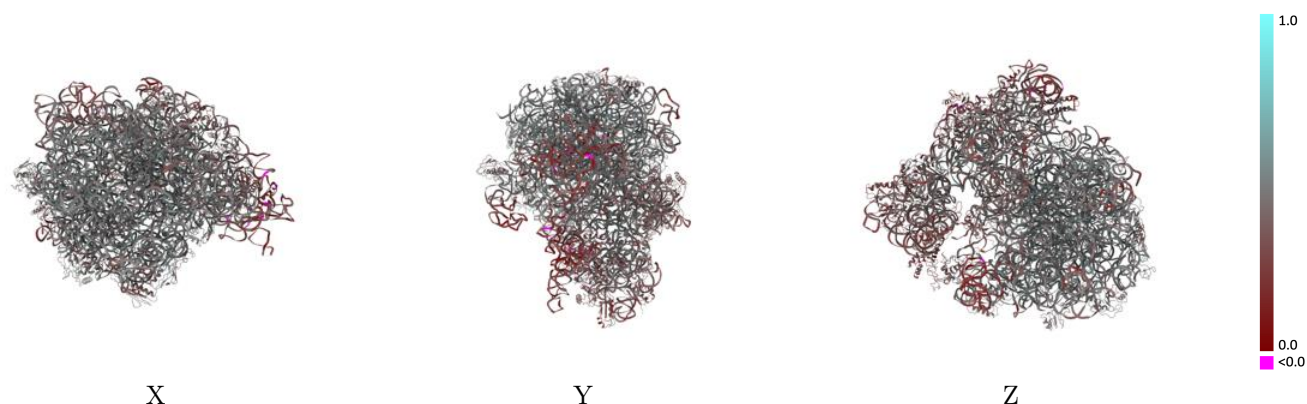
This section contains information regarding the fit between EMDB map EMD-0660 and PDB model 6O90. Per-residue inclusion information can be found in [section 3](#) on [page 13](#).

9.1 Map-model overlay [i](#)



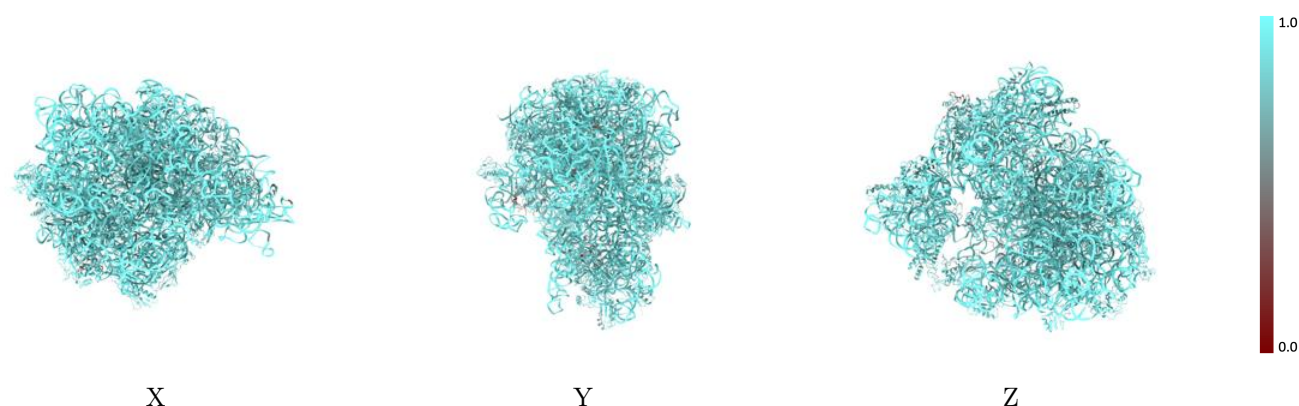
The images above show the 3D surface view of the map at the recommended contour level 0.238 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



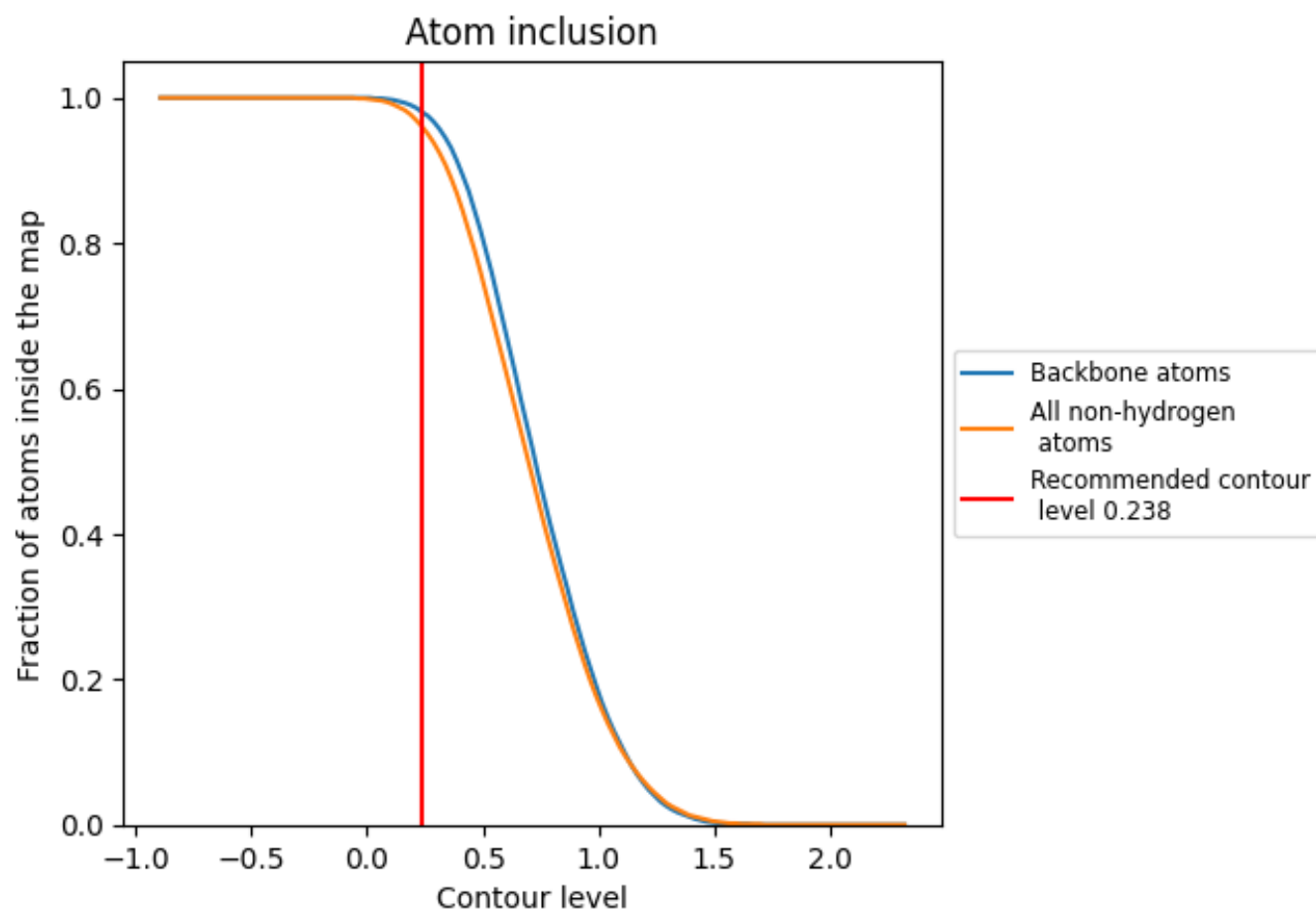
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.238).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 96% 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 (0.238) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9600	 0.4280
0	 0.9180	 0.4660
1	 0.8830	 0.3050
2	 0.9620	 0.4940
3	 0.9060	 0.4700
4	 0.9140	 0.4700
5	 0.9030	 0.4930
6	 0.9520	 0.4820
A	 0.9890	 0.4580
B	 0.9950	 0.4200
C	 0.9300	 0.4780
D	 0.9400	 0.4980
E	 0.9420	 0.4690
F	 0.8760	 0.3670
G	 0.9000	 0.4090
K	 0.9510	 0.4820
L	 0.9170	 0.4930
M	 0.9460	 0.4790
N	 0.8380	 0.4370
O	 0.9360	 0.4670
P	 0.8960	 0.3990
Q	 0.9260	 0.4810
R	 0.9260	 0.4490
S	 0.9160	 0.4630
T	 0.9340	 0.4850
U	 0.9190	 0.4580
V	 0.9000	 0.4450
W	 0.3410	 0.2950
X	 0.9280	 0.4920
Y	 0.9250	 0.4820
Z	 0.9290	 0.4210
a	 0.9910	 0.3940
c	 0.8900	 0.3580
d	 0.7230	 0.2670
e	 0.9010	 0.4220



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Chain	Atom inclusion	Q-score
f	 0.8800	 0.3790
g	 0.8730	 0.3020
h	 0.9220	 0.4100
i	 0.9380	 0.3550
j	 0.8740	 0.3470
k	 0.8780	 0.3970
l	 0.8400	 0.3770
m	 0.8740	 0.2700
n	 0.9430	 0.3970
o	 0.8810	 0.3720
p	 0.9020	 0.3740
q	 0.9030	 0.4190
r	 0.9100	 0.3700
s	 0.8850	 0.3060
t	 0.8160	 0.3340